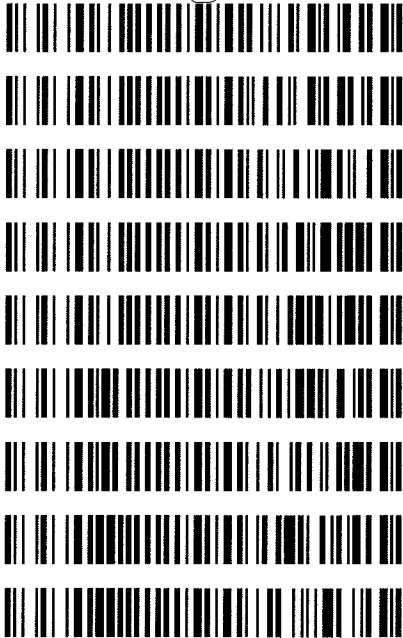


Worklist: 2089

| <u>LAB CASE</u> | <u>ITEM</u> | <u>TASK ID</u> | <u>DESCRIPTION</u> |
|-----------------|-------------|----------------|------------------------------|
| C2017-0502 | 1 | 102768 | AM 28 Blood Multi-Drug Quant |
| C2017-1110 | 1 | 102771 | AM 28 Blood Multi-Drug Quant |
| C2017-2481 | 1 | 103267 | AM 28 Blood Multi-Drug Quant |
| C2017-2531 | 2 | 103266 | AM 28 Blood Multi-Drug Quant |
| C2017-2535 | 1 | 103265 | AM 28 Blood Multi-Drug Quant |
| M2017-1058 | 1 | 102767 | AM 28 Blood Multi-Drug Quant |
| M2017-4285 | 1 | 102192 | AM 28 Blood Multi-Drug Quant |
| P2017-1829 | 4 | 102772 | AM 28 Blood Multi-Drug Quant |
| P2017-2254 | 1 | 102770 | AM 28 Blood Multi-Drug Quant |

BWylee



Multi-Drug Quantitation in Blood by LC-MS/M

Extraction Date: 12-19-17
Plate lot#: 0507680

Analyst: Anne Nord
Plate Expiration: 5/31/2018 6/5/2018 amn

Mobile phase A: 5mM Amm Form + 0.01% FA
0.5M Ammonium Hydroxide
20% Methanol in Water

Mobile phase B: 0.01% Formic Acid in MeOH
Ethyl Acetate

Blank Blood Lot: 321632-1 **Column:** Agilent 120 EC-C18 (2.1x 100-4um)
LCMS-QQQ ID: 62340

Pre-Analytic:

- 1. Check levels of mobile phases and needle wash refill as needed. Ensure waste is not full.
- 2. Ensure correct column is installed and begin mobile phase flow allow to equilibrate ~ 30 minutes.
- 3. Create worklist:

Analytic:

- 1. Remove standards, plate, controls, and samples from cold storage. Allow to reach room temperature.
- 2. Pipette **250µL blood (calibrated pipette) Pipette ID: 2609543** in wells of analytical (standards) plate.
- 3. Place on shaking incubator at ambient temp., 900rpm for 15 minutes. *Shaker ID: 66759*
- 4. Pipette **250µL 00.5M ammonium hydroxide** in wells of analytical plate.
- 5. Place on shaking incubator at ambient temp., 900rpm for 15 minutes.
- 6. Transfer **300µL of blood+base** mixture to corresponding wells of SLE+ plate.
- 7. Apply positive pressure for approx. 10-15 seconds (or until no liquid remains on top of sorbent).
(Load at 85-100 PSI- Selector to the right) Manifold ID: 66792
- 8. Wait 5 minutes.
- 9. Add **900uL ethyl acetate.**
- 10. Wait 5 minutes.
- 11. Apply positive pressure for approx. 15 seconds. **(10-15 PSI- Selector to the left).**
- 12. Add **900uL ethyl acetate.**
- 13. Wait 5 minutes.
- 14. Apply positive pressure for approx. 15 seconds. **(10-15 PSI- Selector to the left).**
- 15. Remove plate containing eluate. Place on SPE Dry and evaporate to dryness at approx. 35°C.
SPE Dry ID: 66819
- 16. Reconstitute in **100µL 20% MeOH** and heat seal plate with foil. Place in autosampler and run worklist.

Post-Analytic

- 1. Create batch and process data.
Worklist path: 121917 mdq p1 and p2 Batch Name: 121917 mdq p2 b
- 2. Make necessary changes to integration limits
- 3. Integration linear and r^2 values ≥ 0.98 for each analyte.
- 4. For unknown samples and controls: response ratio within 20% of average of controls and standards, RT within +/- 5%, S/N for primary and secondary transitions > 10 .
- 5. Did all QCs pass for each analyte? Y/N see comments
- 6. Central File Packet to include: LIMS Worklist, Method Checklist, Calibration and Control Reports

COMMENTS: Protriptyline not evaluated,
100 level calibrator did not inject used 100 ac for 100 level calibrator

Curve ranges: Benzoylcegonine 5-500 andanasetron 5-250
Chlor diazepoxide 5-500 Fluazepam 5-100
Meto prolool 5-500 Phenytoin 5-500
Mirtazapine 5-500

Toxicology AMI method 28 panel 2 external prep information

working solution 10000 ng/ml in meoh mirtazapine, amitriptyline, meperidine

Stock solution 1mg/ml 100 ul each in 9700 ul meOH lot (Fisher 168427)

Ppd 8/7/17 Exp: 2/7/18 lot 2718 by AMN

| Drug | lot | expiration |
|---------------|------------|------------|
| mirtazapine | fe04201503 | 4/1/2020 |
| amitriptyline | fn07081401 | 9/1/2019 |
| meperidine | FE1191502 | 2/1/2020 |

AM 28 control 50 ul working solution lot (2718) in 9950 ul blood lot (321632)

ppd 8/7/17 Exp 11/7/17 by AMN Concentration 50 ng/ml each

ppd 12/5/17 Exp 2/7/18 by AMN Concentration 50 ng/ml each

ISP FORENSICS - Cd'A Instrument # 62340

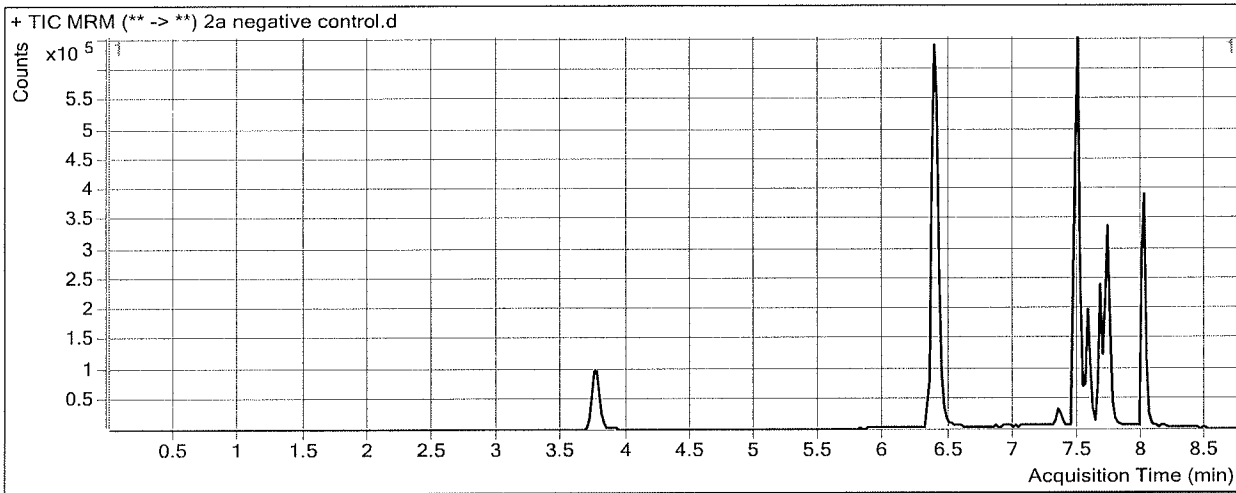
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 08:52 **Data File** 2a negative control.d
Sample Type Sample **Sample Name** 2a negative control
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-E6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------------|
| Nortriptyline | Nortriptyline-D3 | 7.571 | 1397 | 30898 | 0.0452 | 2.7800 <i>LS</i> |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 7356 | 528188 | 0.0139 | 0.9831 <i>LS</i> |

ISP FORENSICS - Cd'A Instrument # 62340

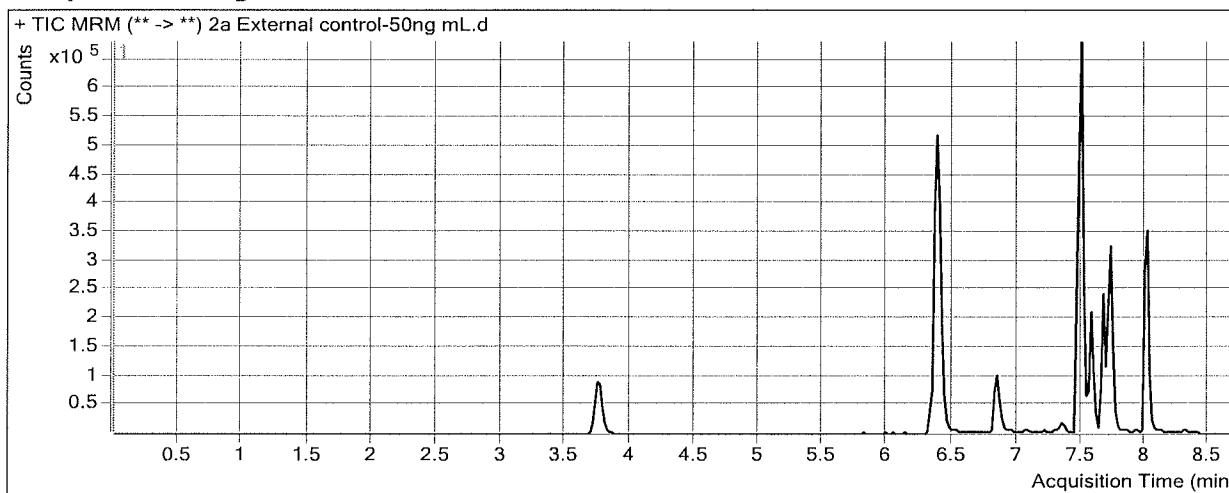
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 09:18 **Data File** 2a External control-50ng mL.d
Sample Type Sample **Sample Name** 2a External control-50ng mL
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-F6 **Sample Info**
Inj Vol -1 **Comment** AM 28 mirtazapine 50ng, amitriptyline 50 ng

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|--------------------|-------|----------|-----------|------------|------------------------------|
| Mirtazapine | Methylphenidate-D4 | 6.845 | 193405 | 680838 | 0.2841 | 31.0084 |
| Maprotiline | Amitriptyline-D3 | 7.567 | 4378 | 16831 | 0.2602 | 16.6816 - ratio out negative |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 28662 | 16831 | 1.7030 | 33.7366 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 511 | 20298 | 0.0252 | 1.9638 (5) |

ISP FORENSICS - Cd'A Instrument # 62340

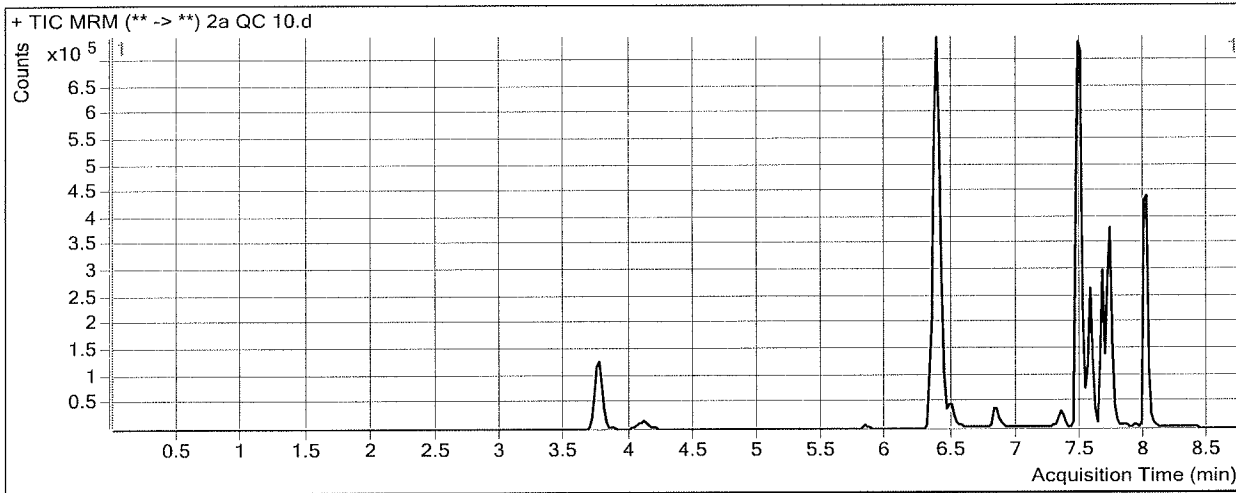
Multi-Drug Confirmatory Analysis Report

| | | | |
|--------------------------|---|----------------------|-----------|
| Batch Data Path | D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin | | |
| Analysis Time | 12/21/2017 10:52 AM | Analyst Name | ISP Tox |
| Report Time | 12/21/2017 10:55 AM | Reporter Name | ISP Tox |
| Last Calib Update | 12/21/2017 10:52 AM | Batch State | Processed |

Analysis Info

| | | | |
|--------------------|------------------|--------------------|--------------------------|
| Acq Time | 2017-12-20 09:43 | Data File | 2a QC 10.d |
| Sample Type | QC | Sample Name | 2a QC 10 |
| Dilution | 1 | Acq Method | MDQ Panel 2a 3-12-17-3.m |
| Position | P1-A6 | Sample Info | |
| Inj Vol | -1 | Comment | AM 28 Panel 2 |

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 97974 | 477671 | 0.2051 | 10.2529 |
| Levamisole | Cocaine-D3 | 4.104 | 51409 | 596649 | 0.0862 | 10.9743 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.831 | 8210 | 7524 | 1.0911 | 11.2116 |
| Zopiclone | Zopiclone-D4 | 6.388 | 20072 | 99300 | 0.2021 | 9.7981 |
| Cocaine | Cocaine-D3 | 6.352 | 140224 | 596649 | 0.2350 | 9.9971 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 205415 | 954922 | 0.2151 | 9.9554 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 32993 | 954922 | 0.0346 | 10.2745 |
| Ondansetron | Zopiclone-D4 | 6.496 | 73934 | 99300 | 0.7446 | 11.1056 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 76531 | 954922 | 0.0801 | 10.4902 |
| Mitragynine | Doxepin-D3 | 7.272 | 3943 | 73020 | 0.0540 | 10.4954 |
| Doxepin | Doxepin-D3 | 7.344 | 13578 | 73020 | 0.1859 | 10.3144 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 329939 | 1762175 | 0.1872 | 10.0189 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 4422 | 24850 | 0.1780 | 10.9419 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 13703 | 24850 | 0.5514 | 10.0324 |
| Estazolam | Estazolam-D5 | 7.563 | 79437 | 375156 | 0.2117 | 10.6290 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 5145 | 24390 | 0.2110 | 9.5394 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 89669 | 540604 | 0.1659 | 10.6533 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 45460 | 457000 | 0.0995 | 10.5796 |
| Midazolam | Midazolam-D4 | 7.721 | 24862 | 381640 | 0.0651 | 9.9201 |
| Phenazepam | Phenazepam-D4 | 7.719 | 6577 | 26515 | 0.2481 | 8.5702 |
| Prazepam | Prazepam-D5 | 8.004 | 209214 | 957544 | 0.2185 | 10.1995 |

ISP FORENSICS - Cd'A Instrument # 62340

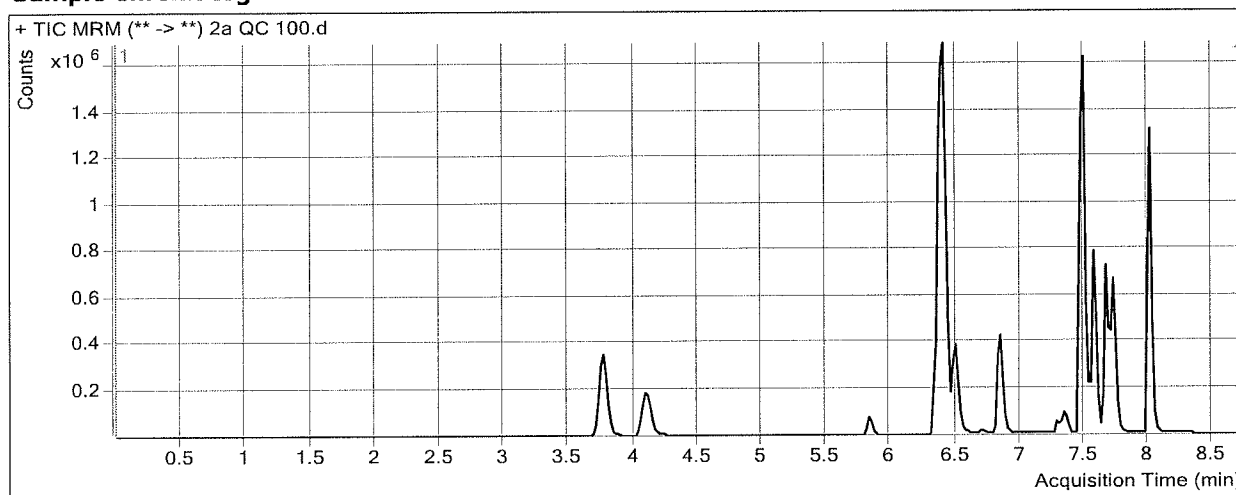
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 13:19 **Data File** 2a QC 100.d
Sample Type Calibration **Sample Name** 2a QC 100
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-B6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|--------------------------|-----------------------------|------------------|------------------|------------------|-------------------|---|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 993530 | 480164 | 2.0691 | 100.7316 |
| Levamisole | Cocaine-D3 | 4.084 | 541586 | 575388 | 0.9413 | 95.1186 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.831 | 107840 | 8039 | 13.4141 | 118.3068 |
| Zopiclone | Zopiclone-D4 | 6.388 | 158480 | 74267 | 2.1339 | 102.1392 |
| Cocaine | Cocaine-D3 | 6.352 | 1385142 | 575388 | 2.4073 | 103.2330 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 2063416 | 958818 | 2.1520 | 98.8545 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 331346 | 958818 | 0.3456 | 109.1428 |
| Ondansetron | Zopiclone-D4 | 6.496 | 754658 | 74267 | 10.1614 | 113.0956 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 834699 | 958818 | 0.8705 | 90.0179 |
| Mitragynine | Doxepin-D3 | 7.272 | 40965 | 73461 | 0.5576 | 95.8049 |
| Doxepin | Doxepin-D3 | 7.344 | 134130 | 73461 | 1.8259 | 98.7124 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 2787106 | 1524226 | 1.8285 | 99.1910 |
| Protriptyline | Protriptyline-D3 | 7.531 | 67598 | 34658 | 1.9504 | 93.4506 <i>not evaluated</i> |
| Maprotiline | Amitriptyline-D3 | 7.547 | 42012 | 27735 | 1.5148 | 104.3019 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 138825 | 27735 | 5.0054 | 101.7184 |
| Estazolam | Estazolam-D5 | 7.563 | 665422 | 310306 | 2.1444 | 92.1486 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 68767 | 26859 | 2.5603 | 105.3476 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 791086 | 468191 | 1.6897 | 107.6338 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 313733 | 368959 | 0.8503 | 106.2799 |
| Midazolam | Midazolam-D4 | 7.721 | 249400 | 359359 | 0.6940 | 98.4015 |
| Phenazepam | Phenazepam-D4 | 7.739 | 43373 | 16726 | 2.5931 | 100.8255 |
| Prazepam | Prazepam-D5 | 8.004 | 1824769 | 852875 | 2.1396 | 105.6085 |

ISP FORENSICS - Cd'A Instrument # 62340

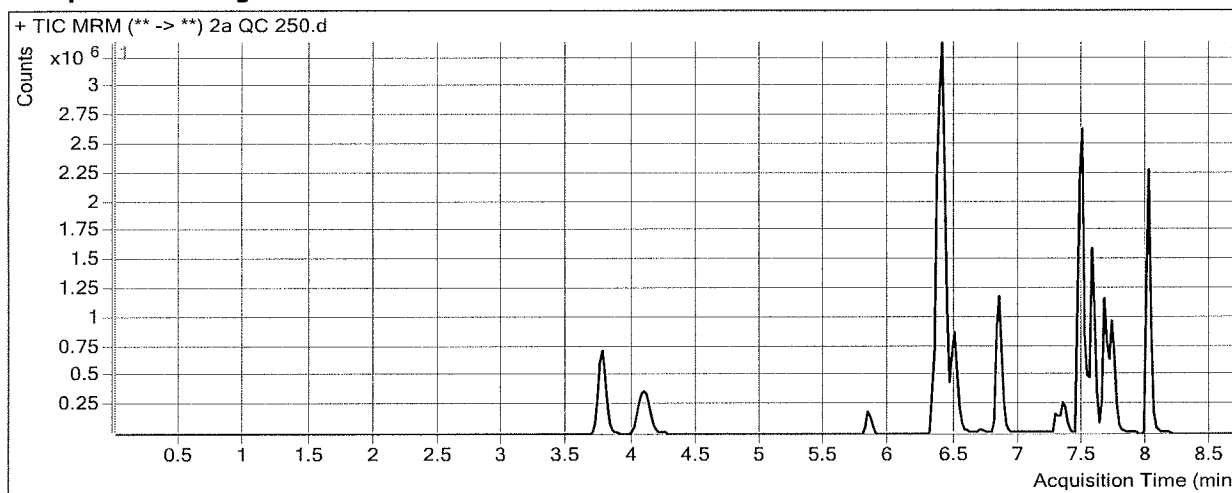
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

| | | | |
|--------------------|------------------|--------------------|--------------------------|
| Acq Time | 2017-12-20 13:45 | Data File | 2a QC 250.d |
| Sample Type | QC | Sample Name | 2a QC 250 |
| Dilution | 1 | Acq Method | MDQ Panel 2a 3-12-17-3.m |
| Position | P1-C6 | Sample Info | |
| Inj Vol | -1 | Comment | AM 28 Panel 2 |

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|--------------------------|-----------------------------|------------------|-------------------|------------------|-------------------|--|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 2417526 | 463814 | 5.2123 | 253.2959 |
| Levamisole | Cocaine-D3 | 4.084 | 1336317 | 555761 | 2.4045 | 239.1061 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.851 | 286896 | 7029 | 40.8141 | 356.4307 |
| Zopiclone | Zopiclone-D4 | 6.388 | 310779 | 61853 | 5.0245 | 240.3128 |
| Cocaine | Cocaine-D3 | 6.352 | 3278167 | 555761 | 5.8985 | 253.0763 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 5203974 | 969858 | 5.3657 | 246.3521 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 727583 | 969858 | 0.7502 | 237.7613 |
| Ondansetron | Zopiclone-D4 | 6.496 | 1775598 | 61853 | 28.7068 | 313.9543 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 2354226 | 969858 | 2.4274 | 246.6619 |
| Mitragynine | Doxepin-D3 | 7.272 | 143846 | 99134 | 1.4510 | 247.1341 |
| Doxepin | Doxepin-D3 | 7.344 | 512234 | 99134 | 5.1671 | 278.8174 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 5583110 | 1214609 | 4.5966 | 249.5815 |
| Protriptyline | Protriptyline-D3 | 7.531 | 244498 | 45231 | 5.4055 | 256.4443 <i>Not evaluated</i> |
| Maprotiline | Amitriptyline-D3 | 7.547 | 152591 | 40157 | 3.7998 | 263.8881 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 512274 | 40157 | 12.7567 | 261.2822 |
| Estazolam | Estazolam-D5 | 7.563 | 1313900 | 234980 | 5.5915 | 237.5483 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 227833 | 39068 | 5.8318 | 238.7590 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 1596826 | 382232 | 4.1776 | 265.9782 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 548597 | 265663 | 2.0650 | 261.1006 |
| Midazolam | Midazolam-D4 | 7.721 | 570555 | 325706 | 1.7517 | 247.2243 |
| Phenazepam | Phenazepam-D4 | 7.739 | 69650 | 10702 | 6.5079 | 254.8336 |
| Prazepam | Prazepam-D5 | 8.004 | 3885118 | 753266 | 5.1577 | 255.5039 |

[Handwritten signature]

ISP FORENSICS - Cd'A Instrument # 62340

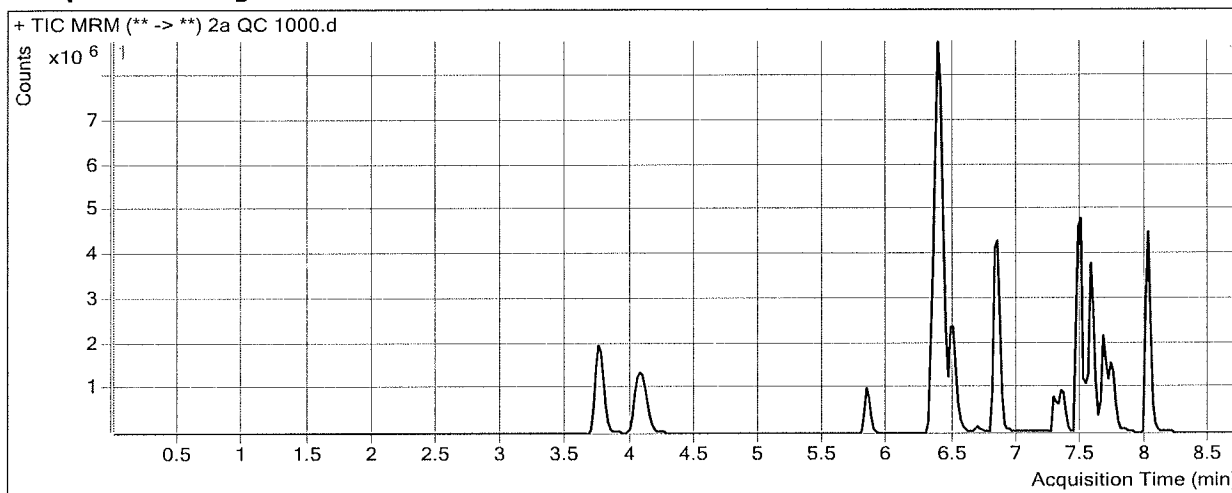
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 14:10 **Data File** 2a QC 1000.d
Sample Type QC **Sample Name** 2a QC 1000
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-D6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|-----------------------------|--------------------------------|------------------|--------------------|-------------------|---------------------|--|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.750 | 8266681 | 395826 | 20.8846 | 1014.0180 |
| Levamisole | Cocaine-D3 | 4.064 | 5031258 | 443713 | 11.3390 | 1118.2988 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.831 | 1529721 | 7113 | 215.0603 | 1870.7446 * |
| Zopiclone | Zopiclone-D4 | 6.368 | 458700 | 22533 | 20.3568 | 973.2160 |
| Cocaine | Cocaine-D3 | 6.352 | 10537203 | 443713 | 23.7478 | 1019.1737 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 15630412 | 704413 | 22.1893 | 1018.5014 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 1776310 | 704413 | 2.5217 | 800.8763 * |
| Ondansetron | Zopiclone-D4 | 6.476 | 5812594 | 22533 | 257.9588 | 2796.8978 * |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 9468001 | 704413 | 13.4410 | 1354.8075 * |
| Mitragynine | Doxepin-D3 | 7.272 | 685444 | 114019 | 6.0117 | 1019.6578 |
| Doxepin | Doxepin-D3 | 7.364 | 2223605 | 114019 | 19.5021 | 1051.5378 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 11744545 | 622964 | 18.8527 | 1024.1134 |
| Protriptyline | Protriptyline-D3 | 7.531 | 834056 | 39440 | 21.1473 | 999.0529 <i>not evaluated</i> * |
| Maprotiline | Amitriptyline-D3 | 7.567 | 549204 | 46319 | 11.8569 | 826.5813 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 2149717 | 46319 | 46.4108 | 954.0625 |
| Estazolam | Estazolam-D5 | 7.563 | 2627450 | 107047 | 24.5448 | 1036.9955 |
| Nortriptyline | Nortriptyline-D3 | 7.591 | 861412 | 34649 | 24.8612 | 1014.7902 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 3545859 | 240724 | 14.7300 | 937.5692 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 964291 | 126707 | 7.6104 | 967.8995 * |
| Midazolam | Midazolam-D4 | 7.721 | 1587752 | 219666 | 7.2280 | 1017.7358 |
| Phenazepam | Phenazepam-D4 | 7.739 | 120409 | 5280 | 22.8040 | 895.9292 |
| Prazepam | Prazepam-D5 | 8.004 | 10063453 | 489763 | 20.5476 | 1019.8387 |

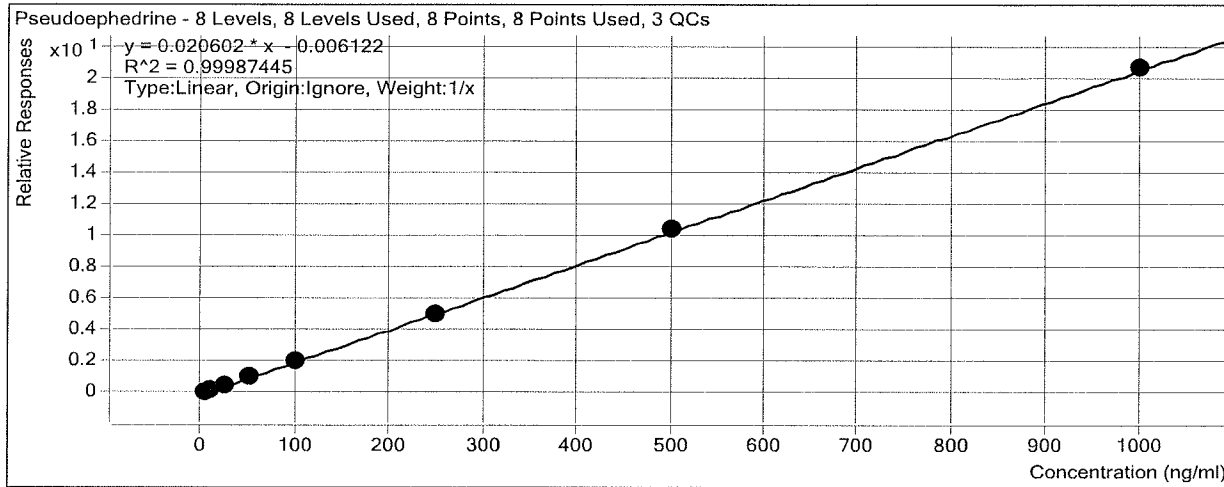
* outside range of curve *

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Pseudoephedrine*
Internal Standard *Pseudoephedrine-D3*



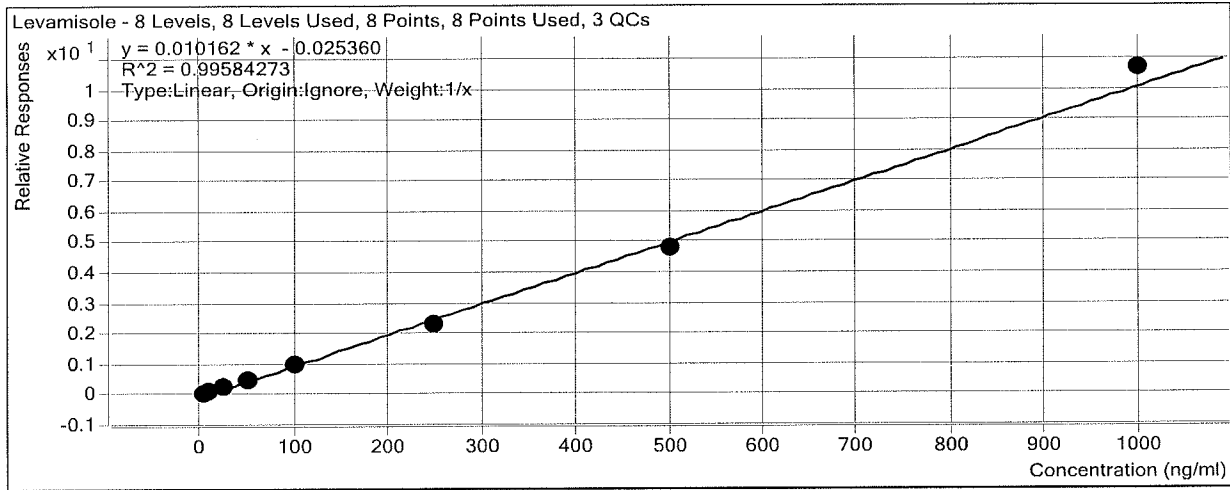
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 102.4 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.0 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 102.5 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.4 | 97.5 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 50.4 | 100.8 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 100.7 | 100.7 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 243.8 | 97.5 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 253.3 | 101.3 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 504.5 | 100.9 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1001.1 | 100.1 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1014.0 | 101.4 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Levamisole
Internal Standard Cocaine-D3



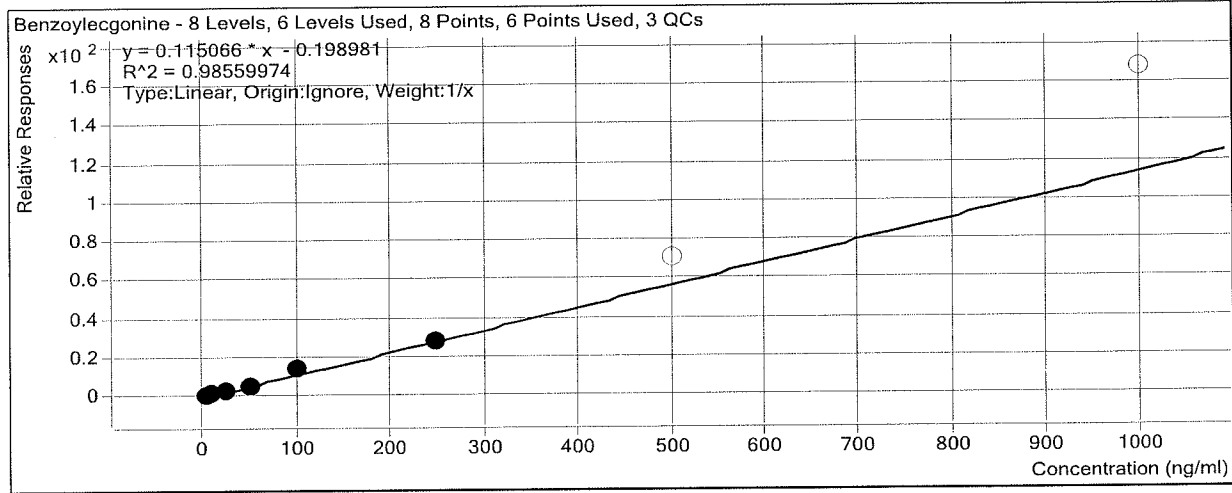
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.3 | 125.6 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 105.4 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 11.0 | 109.7 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.0 | 92.0 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 45.0 | 90.0 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 95.1 | 95.1 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 226.5 | 90.6 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 239.1 | 95.6 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 478.6 | 95.7 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1054.9 | 105.5 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1118.3 | 111.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Benzoyllecgonine*
Internal Standard *Benzoyllecgonine-D8*



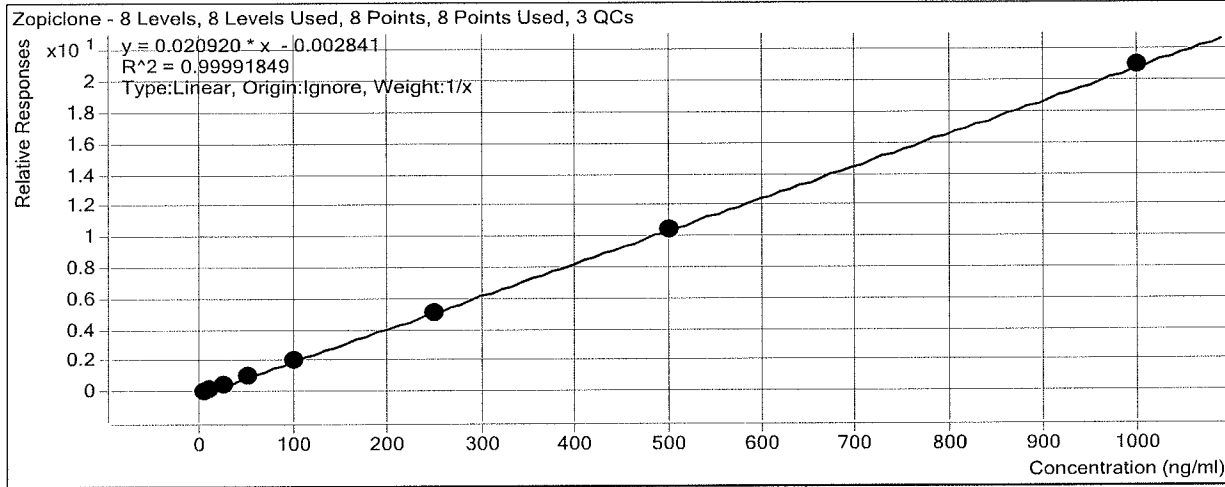
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.5 | 110.5 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 101.8 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 11.2 | 112.1 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 21.3 | 85.1 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 44.0 | 88.1 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 118.3 | 118.3 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 240.7 | 96.3 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 356.4 | 142.6 |
| 2a cal 7-500ng | 7 | <input type="checkbox"/> | 500 | 614.7 | 122.9 |
| 2a cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1460.5 | 146.0 |
| 2a QC 1000 | 8 | <input type="checkbox"/> | 1000 | 1870.7 | 187.1 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Zopiclone
Internal Standard Zopiclone-D4



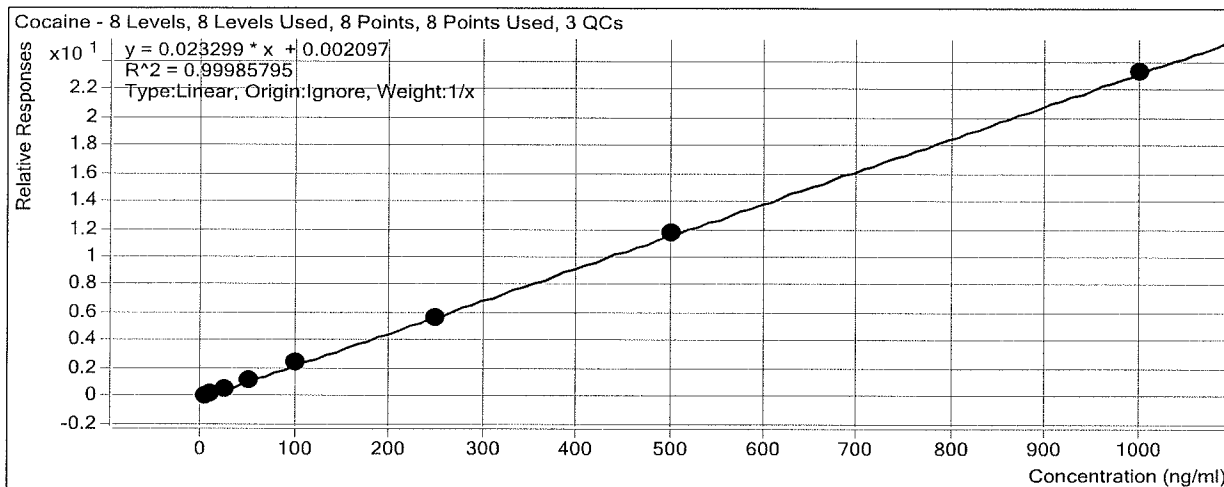
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 102.2 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.2 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.0 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.4 | 97.6 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 50.7 | 101.3 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 102.1 | 102.1 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 246.0 | 98.4 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 240.3 | 96.1 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 499.6 | 99.9 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1002.3 | 100.2 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 973.2 | 97.3 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Cocaine
Internal Standard Cocaine-D3



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.8 | 95.8 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 99.4 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.0 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.1 | 100.5 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.4 | 102.8 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 103.2 | 103.2 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 245.3 | 98.1 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 253.1 | 101.2 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 502.0 | 100.4 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 998.2 | 99.8 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1019.2 | 101.9 |

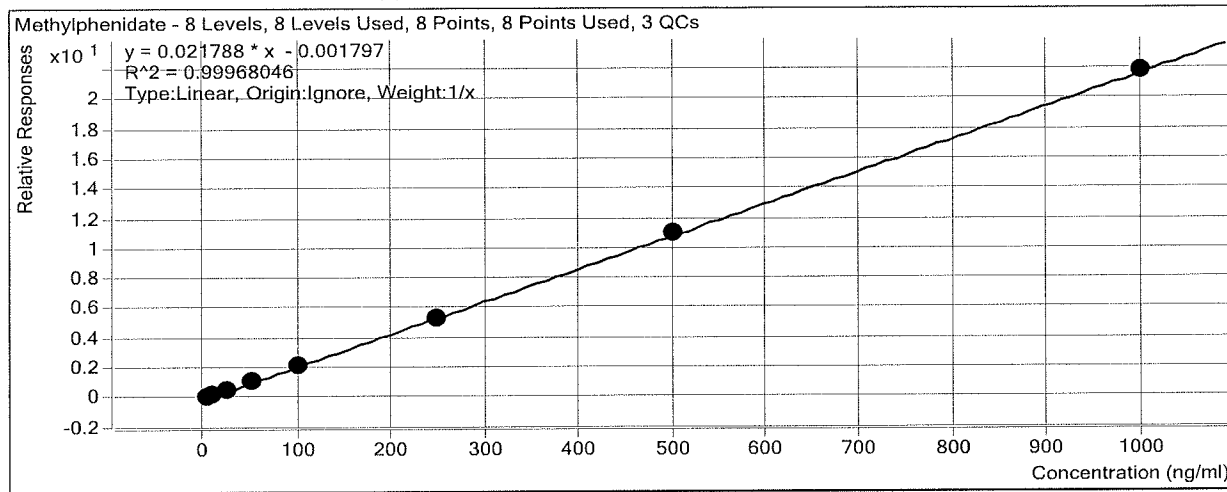
ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM

Analyst Name ISP TOX

Target Compound *Methylphenidate*
Internal Standard *Methylphenidate-D4*



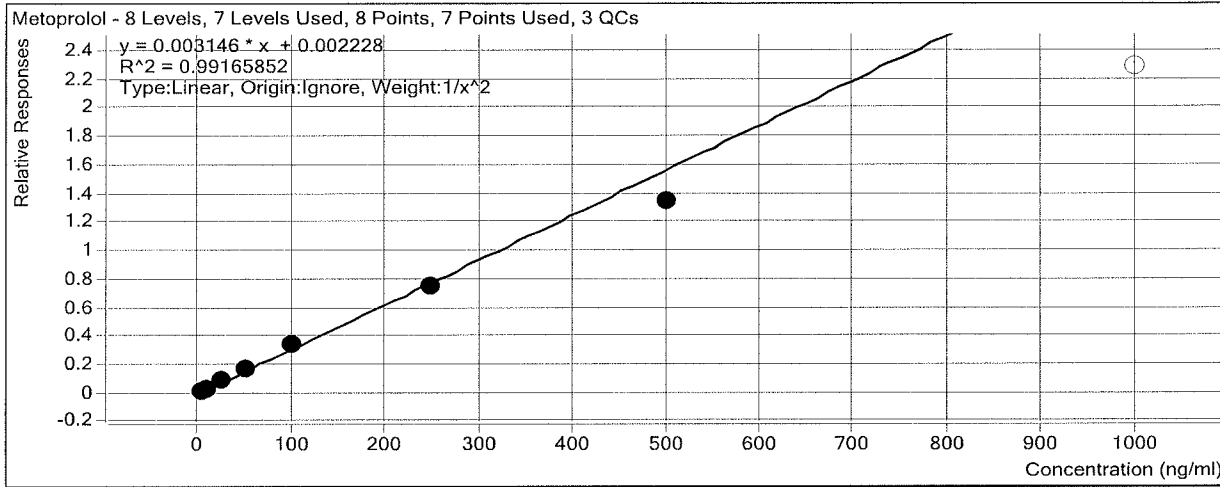
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 103.8 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.0 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 99.6 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.7 | 98.6 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 50.2 | 100.5 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 98.9 | 98.9 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 240.6 | 96.2 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 246.4 | 98.5 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 509.6 | 101.9 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1000.9 | 100.1 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1018.5 | 101.9 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Metoprolol*
Internal Standard *Methylphenidate-D4*



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.9 | 97.3 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 102.5 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 102.7 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.7 | 102.8 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 53.2 | 106.3 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 109.1 | 109.1 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 240.1 | 96.0 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 237.8 | 95.1 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 429.3 | 85.9 |
| 2a cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 725.2 | 72.5 |
| 2a QC 1000 | 8 | <input type="checkbox"/> | 1000 | 800.9 | 80.1 |

ISP Forensics Calibration Curve Report

Batch Data Path

D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update

12/21/2017 10:52 AM

Analyst Name

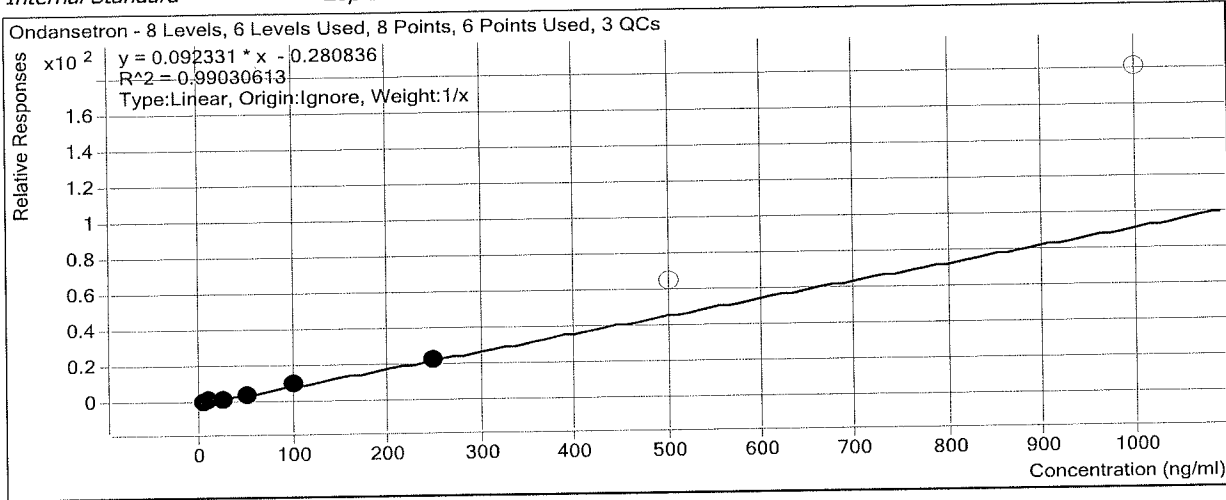
ISP TOX

Target Compound

Ondansetron

Internal Standard

Zopiclone-D4



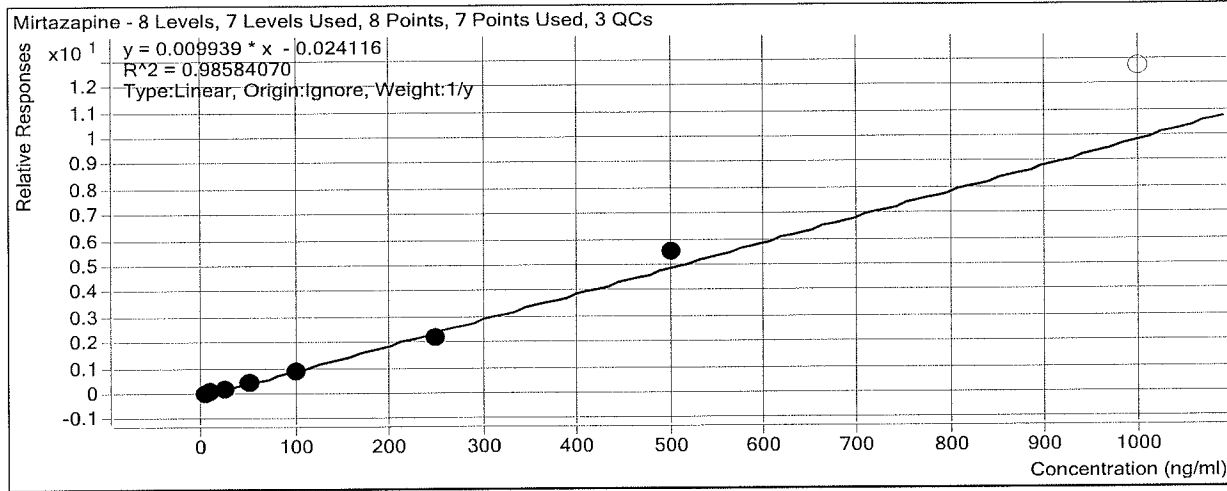
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.2 | 123.1 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.3 | 92.9 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 11.1 | 111.1 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 20.2 | 80.8 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 46.0 | 92.0 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 113.1 | 113.1 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 245.3 | 98.1 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 314.0 | 125.6 |
| 2a cal 7-500ng | 7 | <input type="checkbox"/> | 500 | 698.5 | 139.7 |
| 2a cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1959.1 | 195.9 |
| 2a QC 1000 | 8 | <input type="checkbox"/> | 1000 | 2796.9 | 279.7 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Mirtazapine*
Internal Standard *Methylphenidate-D4*



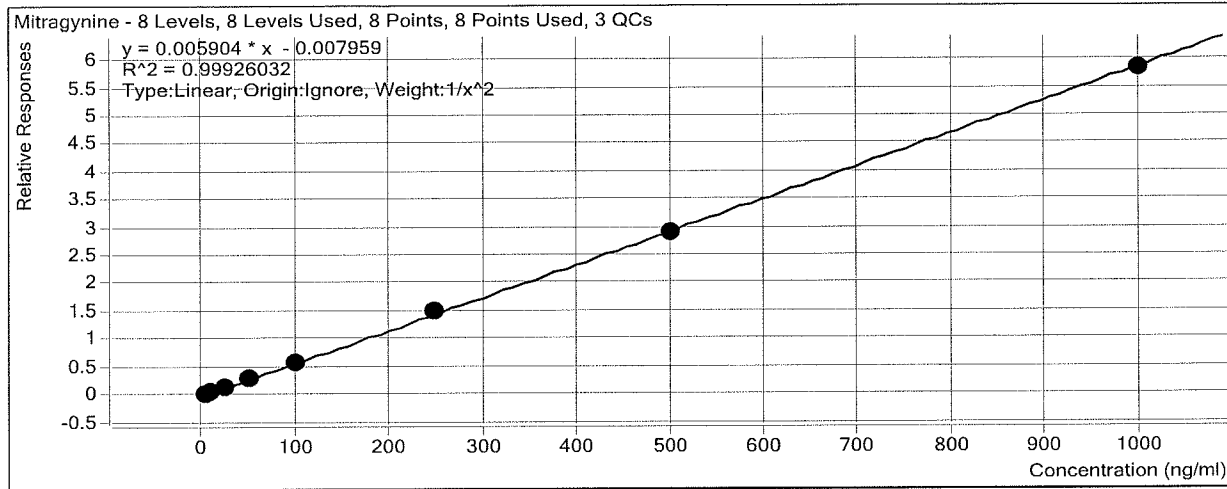
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.3 | 126.9 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.0 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 104.9 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 22.3 | 89.4 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 45.5 | 91.0 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 90.0 | 90.0 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 221.5 | 88.6 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 246.7 | 98.7 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 555.8 | 111.2 |
| 2a cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1278.4 | 127.8 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1354.8 | 135.5 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Mitragynine*
Internal Standard *Doxepin-D3*



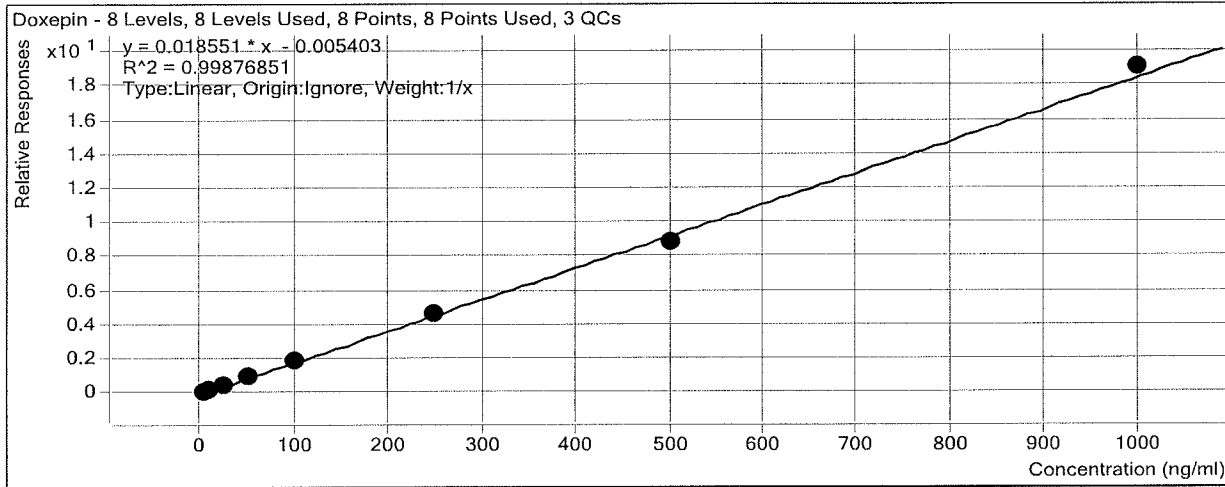
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.0 | 99.0 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.1 | 100.6 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 105.0 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.7 | 102.9 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.6 | 103.1 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 95.8 | 95.8 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 251.2 | 100.5 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 247.1 | 98.9 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 496.2 | 99.2 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 988.6 | 98.9 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1019.7 | 102.0 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Doxepin
Internal Standard Doxepin-D3



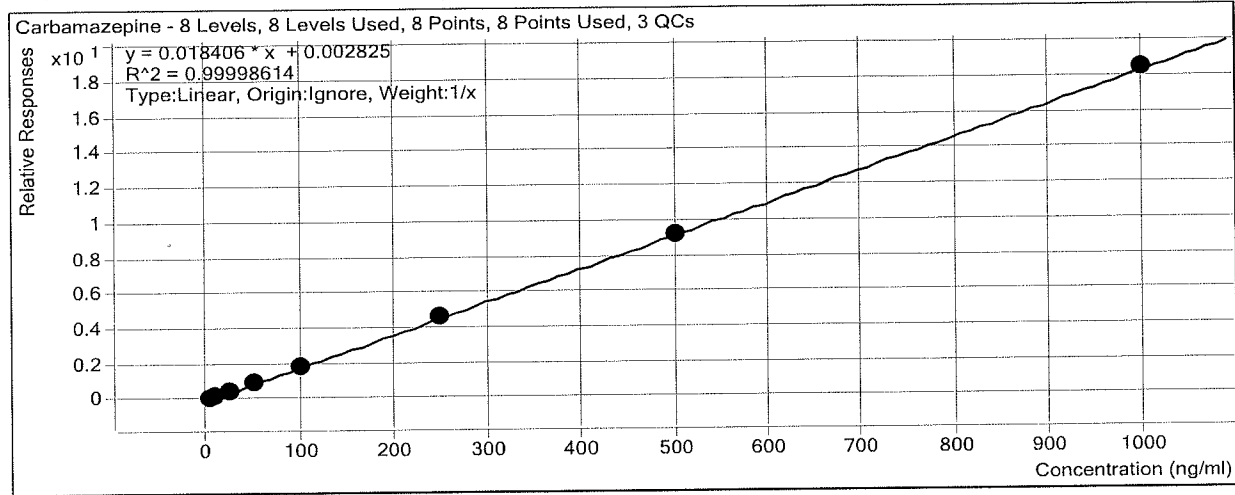
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.0 | 100.7 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 97.6 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.1 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.6 | 102.3 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.9 | 103.8 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 98.7 | 98.7 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 248.8 | 99.5 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 278.8 | 111.5 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 473.8 | 94.8 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1026.4 | 102.6 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1051.5 | 105.2 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Carbamazepine
Internal Standard Carbamazepine-13C6



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 101.4 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 99.5 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.2 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.1 | 100.2 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 49.9 | 99.9 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 99.2 | 99.2 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 249.9 | 100.0 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 249.6 | 99.8 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 498.0 | 99.6 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1002.9 | 100.3 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1024.1 | 102.4 |

ISP Forensics Calibration Curve Report

Batch Data Path

D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update

12/21/2017 10:52 AM

Analyst Name

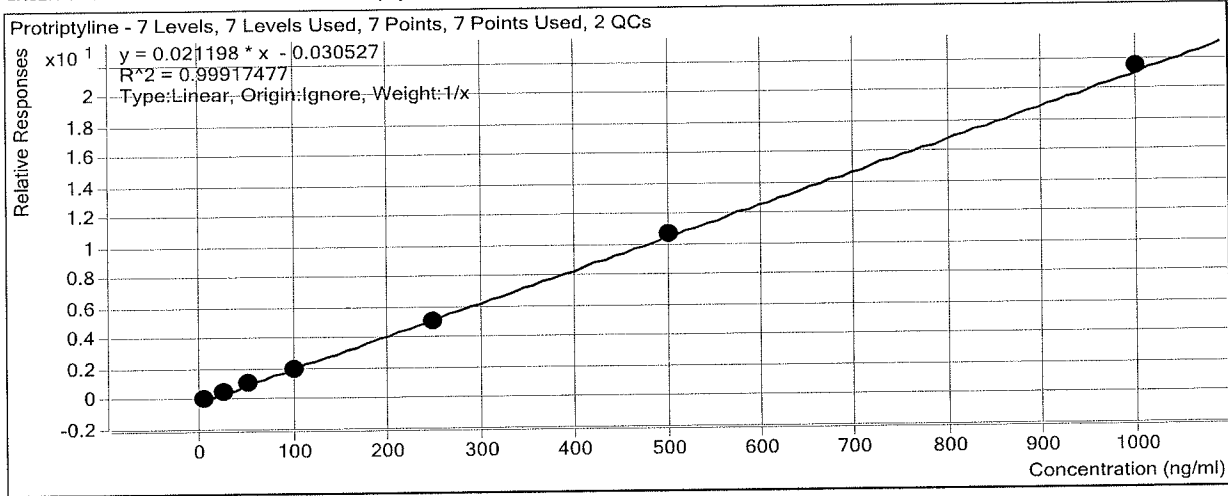
ISP TOX

Target Compound

Protriptyline

Internal Standard

Protriptyline-D3



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.9 | 118.3 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 0.0 | 0.0 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 0.0 | 0.0 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.0 | 91.9 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 48.9 | 97.8 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 93.5 | 93.5 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 239.7 | 95.9 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 256.4 | 102.6 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 506.8 | 101.4 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1012.2 | 101.2 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 999.1 | 99.9 |

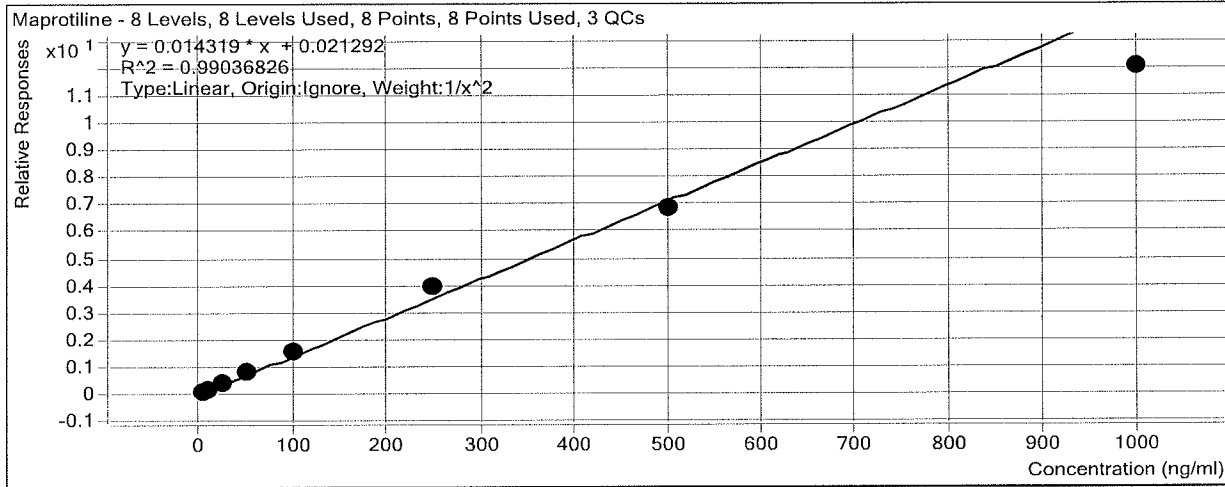
Not evaluated

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Maprotiline*
Internal Standard *Amitriptyline-D3*



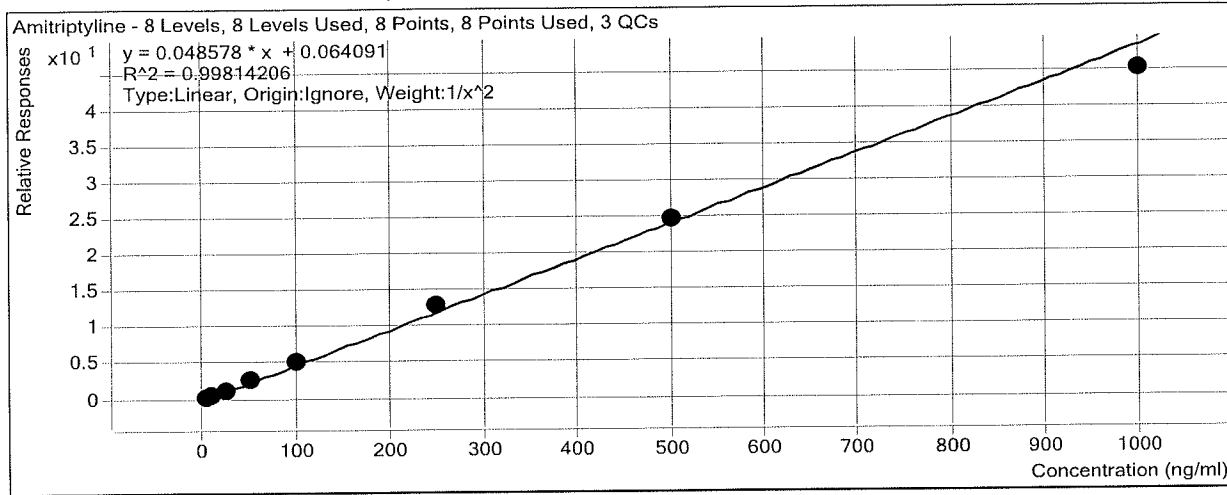
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.9 | 98.2 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 102.9 |
| 2a QC 10 | 2 | <input type="checkbox"/> | 10 | 10.9 | 109.4 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.0 | 95.8 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 54.2 | 108.4 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 104.3 | 104.3 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 277.5 | 111.0 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 263.9 | 105.6 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 476.0 | 95.2 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 841.5 | 84.2 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 826.6 | 82.7 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Amitriptyline
Internal Standard Amitriptyline-D3



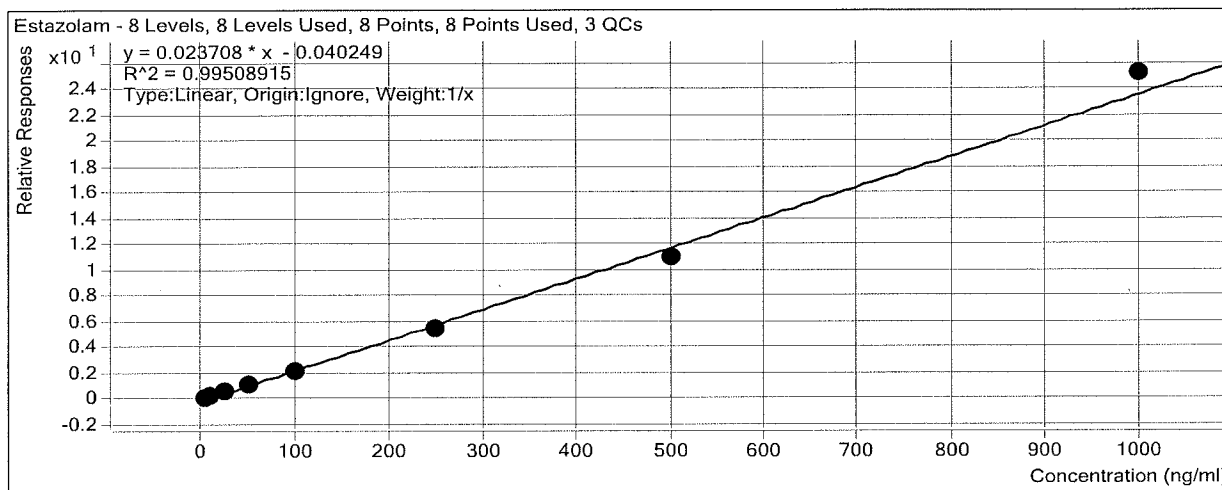
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.0 | 100.8 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.5 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.3 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.5 | 98.2 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 50.6 | 101.1 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 101.7 | 101.7 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 263.7 | 105.5 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 261.3 | 104.5 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 507.9 | 101.6 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 925.9 | 92.6 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 954.1 | 95.4 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Estazolam*
Internal Standard *Estazolam-D5*



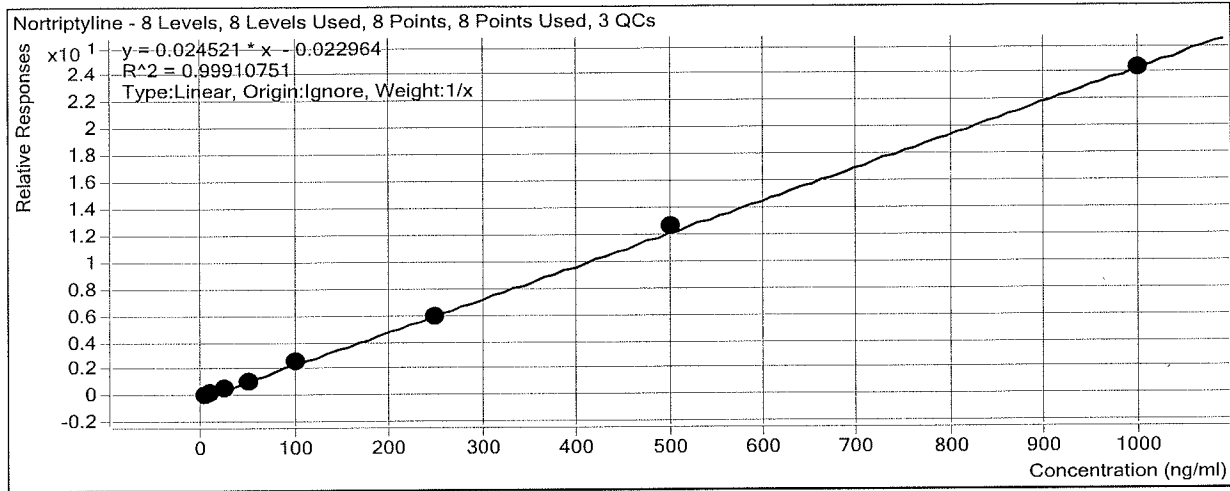
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.1 | 122.0 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.0 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.6 | 106.3 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.5 | 94.0 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 48.7 | 97.5 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 92.1 | 92.1 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 229.0 | 91.6 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 237.5 | 95.0 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 467.8 | 93.6 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1062.5 | 106.2 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1037.0 | 103.7 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Nortriptyline*
Internal Standard *Nortriptyline-D3*



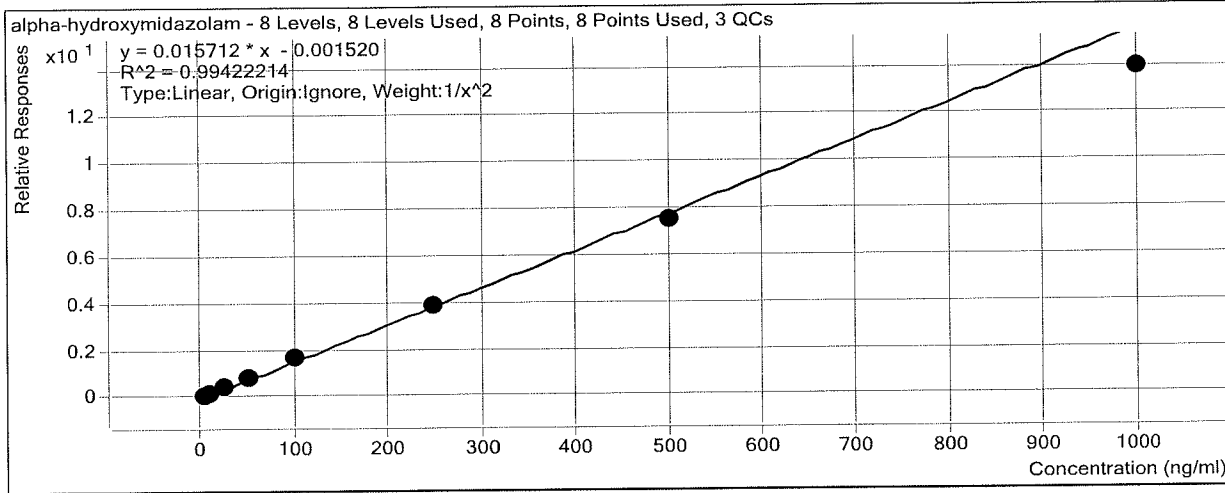
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.9 | 118.4 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.4 | 93.6 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.5 | 95.4 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 22.3 | 89.2 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 47.1 | 94.2 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 105.3 | 105.3 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 242.6 | 97.0 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 238.8 | 95.5 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 513.6 | 102.7 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 993.8 | 99.4 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1014.8 | 101.5 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *alpha-hydroxymidazolam*
Internal Standard *alpha-hydroxymidazolam-D4*



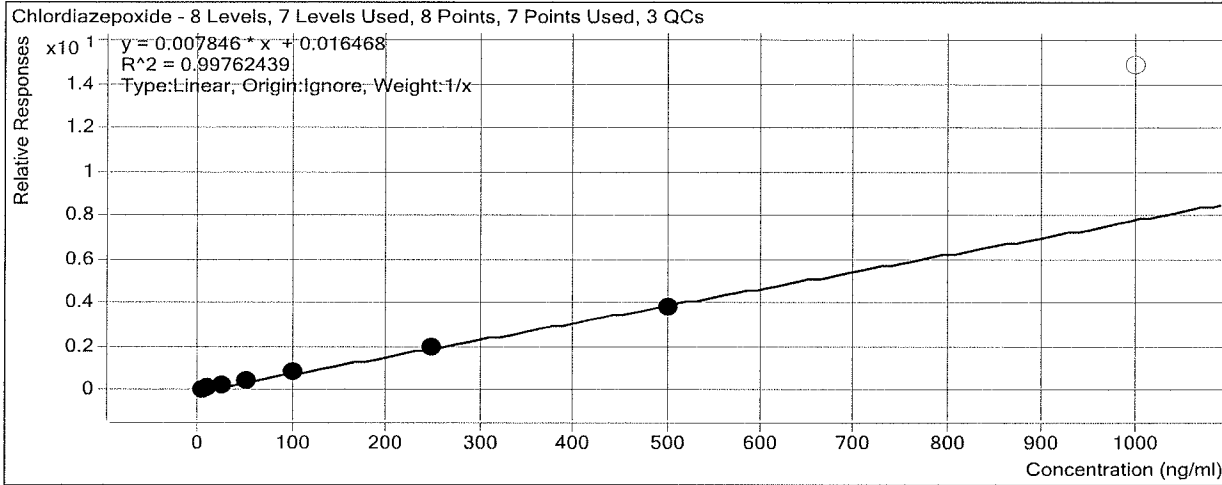
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.8 | 95.6 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.6 | 105.9 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.7 | 106.5 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.8 | 103.1 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 52.7 | 105.5 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 107.6 | 107.6 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 245.9 | 98.4 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 266.0 | 106.4 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 476.4 | 95.3 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 886.1 | 88.6 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 937.6 | 93.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Chlordiazepoxide*
Internal Standard *Chlordiazepoxide-D5*



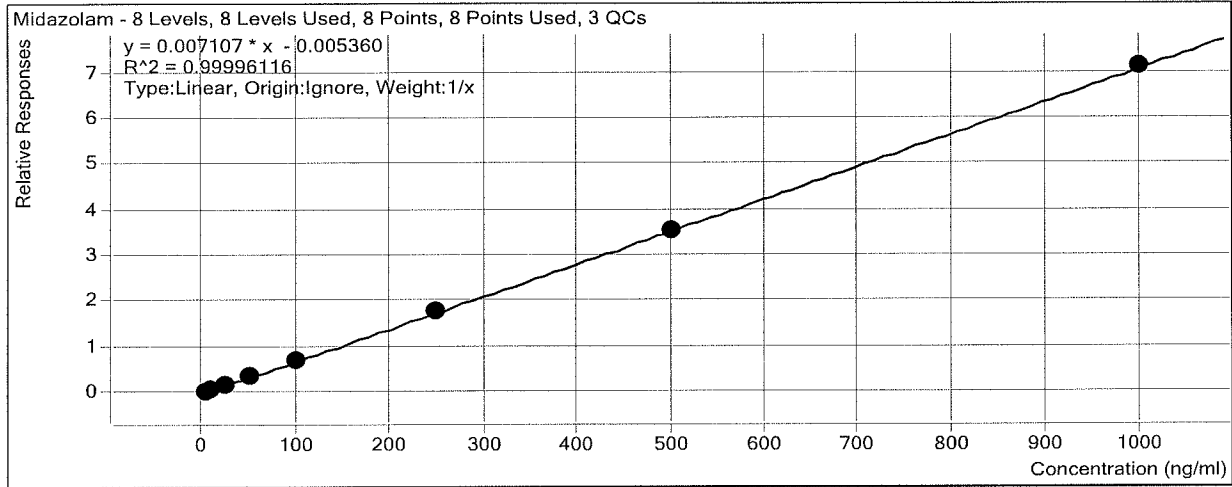
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.1 | 81.7 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.4 | 93.5 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.6 | 105.8 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 27.4 | 109.5 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 55.5 | 111.0 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 106.3 | 106.3 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 252.6 | 101.1 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 261.1 | 104.4 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 484.8 | 97.0 |
| 2a cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1890.7 | 189.1 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 967.9 | 96.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound *Midazolam*
Internal Standard *Midazolam-D4*



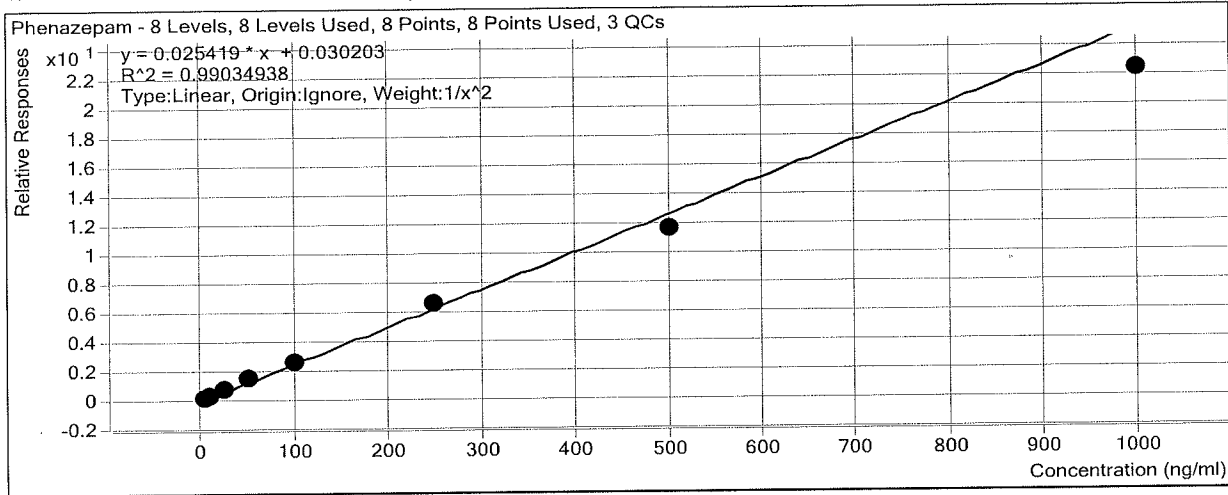
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.3 | 105.0 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 97.9 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 99.2 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.0 | 100.0 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 49.3 | 98.5 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 98.4 | 98.4 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 249.1 | 99.6 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 247.2 | 98.9 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 502.3 | 100.5 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1000.9 | 100.1 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1017.7 | 101.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Phenazepam
Internal Standard Phenazepam-D4



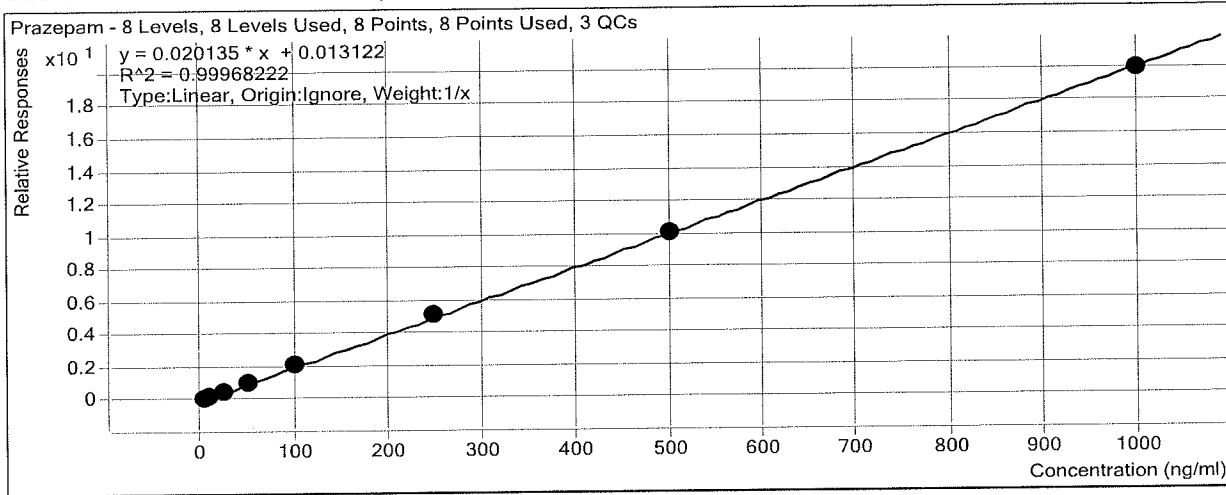
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.9 | 98.6 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.7 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 8.6 | 85.7 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.6 | 102.3 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 58.4 | 116.8 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 100.8 | 100.8 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 257.9 | 103.2 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 254.8 | 101.9 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 458.6 | 91.7 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 880.1 | 88.0 |
| 2a QC 1000 | 8 | <input type="checkbox"/> | 1000 | 895.9 | 89.6 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin

Last Calib Update 12/21/2017 10:52 AM **Analyst Name** ISP TOX

Target Compound Prazepam
Internal Standard Prazepam-D5



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2a cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.6 | 91.7 |
| 2a cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.0 |
| 2a QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.0 |
| 2a cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.2 | 101.0 |
| 2a cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.7 | 103.4 |
| 2a QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 105.6 | 105.6 |
| 2a cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 252.9 | 101.2 |
| 2a QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 255.5 | 102.2 |
| 2a cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 500.9 | 100.2 |
| 2a cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 989.3 | 98.9 |
| 2a QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1019.8 | 102.0 |

ISP FORENSICS - Cd'A Instrument # 62340

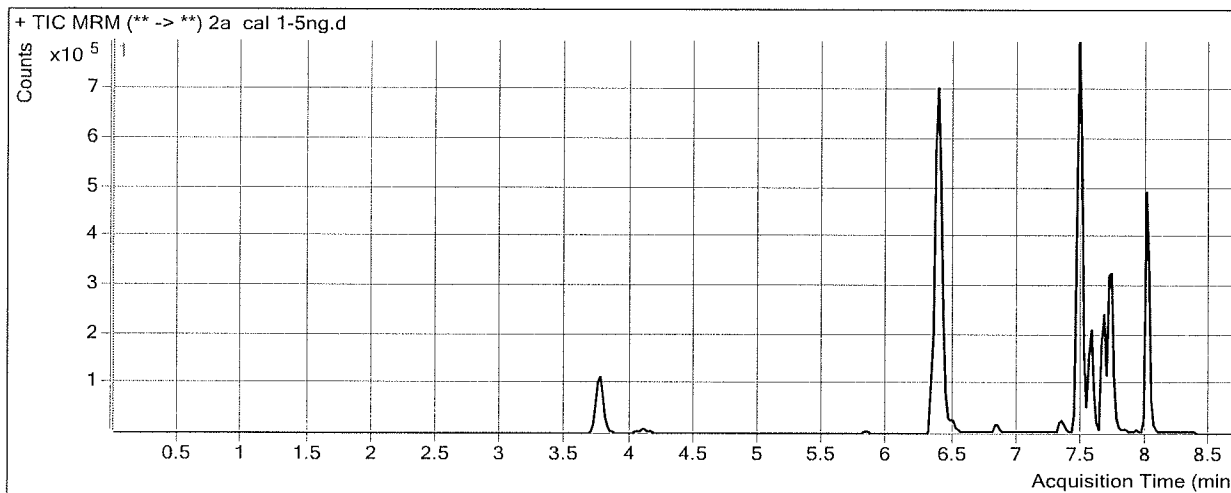
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 05:16 **Data File** 2a cal 1-5ng.d
Sample Type Calibration **Sample Name** 2a cal 1-5ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-A5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 47327 | 476412 | 0.0993 | 5.1191 |
| Levamisole | Cocaine-D3 | 4.084 | 24539 | 638387 | 0.0384 | 6.2781 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.831 | 3433 | 7862 | 0.4366 | 5.5238 |
| Zopiclone | Zopiclone-D4 | 6.388 | 13192 | 126768 | 0.1041 | 5.1104 |
| Cocaine | Cocaine-D3 | 6.352 | 72575 | 638387 | 0.1137 | 4.7894 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 104730 | 941382 | 0.1113 | 5.1886 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 16506 | 941382 | 0.0175 | 4.8653 |
| Ondansetron | Zopiclone-D4 | 6.496 | 36437 | 126768 | 0.2874 | 6.1547 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 36648 | 941382 | 0.0389 | 6.3434 |
| Mitragynine | Doxepin-D3 | 7.272 | 1443 | 67835 | 0.0213 | 4.9513 |
| Doxepin | Doxepin-D3 | 7.344 | 5967 | 67835 | 0.0880 | 5.0330 |
| Carbamazepine | Carbamazepine-13C6 | 7.473 | 169674 | 1764863 | 0.0961 | 5.0698 |
| Protriptyline | Protriptyline-D3 | 7.531 | 2956 | 31151 | 0.0949 | 5.9170 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 2040 | 22270 | 0.0916 | 4.9115 |
| Amitriptyline | Amitriptyline-D3 | 7.547 | 6882 | 22270 | 0.3090 | 5.0421 |
| Estazolam | Estazolam-D5 | 7.563 | 40899 | 391975 | 0.1043 | 6.0988 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 2799 | 22890 | 0.1223 | 5.9224 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 40506 | 550259 | 0.0736 | 4.7817 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 22499 | 463800 | 0.0485 | 4.0840 |
| Midazolam | Midazolam-D4 | 7.721 | 12375 | 387253 | 0.0320 | 5.2503 |
| Phenazepam | Phenazepam-D4 | 7.719 | 4049 | 26041 | 0.1555 | 4.9285 |
| Prazepam | Prazepam-D5 | 8.004 | 105644 | 1001836 | 0.1055 | 4.5855 |

ISP FORENSICS - Cd'A Instrument # 62340

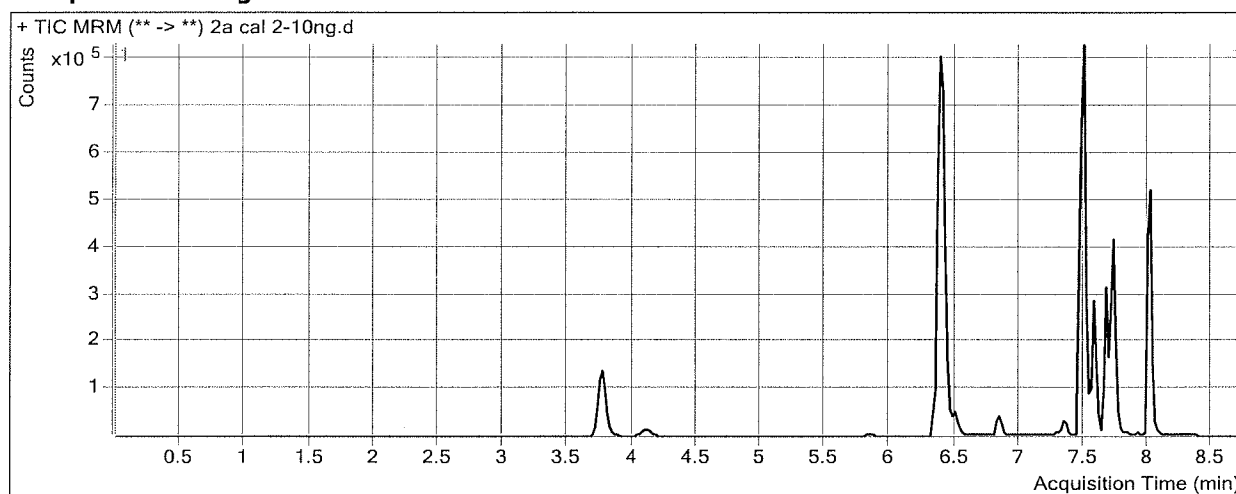
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 05:41 **Data File** 2a cal 2-10ng.d
Sample Type Calibration **Sample Name** 2a cal 2-10ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-B5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 103042 | 515536 | 0.1999 | 9.9988 |
| Levamisole | Cocaine-D3 | 4.104 | 54742 | 669476 | 0.0818 | 10.5418 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.851 | 7981 | 8212 | 0.9719 | 10.1762 |
| Zopiclone | Zopiclone-D4 | 6.388 | 27777 | 137097 | 0.2026 | 9.8208 |
| Cocaine | Cocaine-D3 | 6.352 | 156481 | 669476 | 0.2337 | 9.9421 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 229293 | 1060699 | 0.2162 | 10.0041 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 36578 | 1060699 | 0.0345 | 10.2538 |
| Ondansetron | Zopiclone-D4 | 6.496 | 79123 | 137097 | 0.5771 | 9.2923 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 82971 | 1060699 | 0.0782 | 10.2970 |
| Mitragynine | Doxepin-D3 | 7.272 | 3899 | 75818 | 0.0514 | 10.0603 |
| Doxepin | Doxepin-D3 | 7.344 | 13322 | 75818 | 0.1757 | 9.7631 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 350262 | 1884302 | 0.1859 | 9.9456 |
| Maprotiline | Amitriptyline-D3 | 7.567 | 3909 | 23179 | 0.1687 | 10.2921 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 12573 | 23179 | 0.5424 | 9.8467 |
| Estazolam | Estazolam-D5 | 7.563 | 84379 | 413629 | 0.2040 | 10.3023 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 5174 | 25037 | 0.2066 | 9.3633 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 94221 | 571420 | 0.1649 | 10.5909 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.723 | 44603 | 496465 | 0.0898 | 9.3520 |
| Midazolam | Midazolam-D4 | 7.721 | 26479 | 412367 | 0.0642 | 9.7889 |
| Phenazepam | Phenazepam-D4 | 7.739 | 8557 | 30450 | 0.2810 | 9.8675 |
| Prazepam | Prazepam-D5 | 8.004 | 226046 | 1073746 | 0.2105 | 9.8038 |

ISP FORENSICS - Cd'A Instrument # 62340

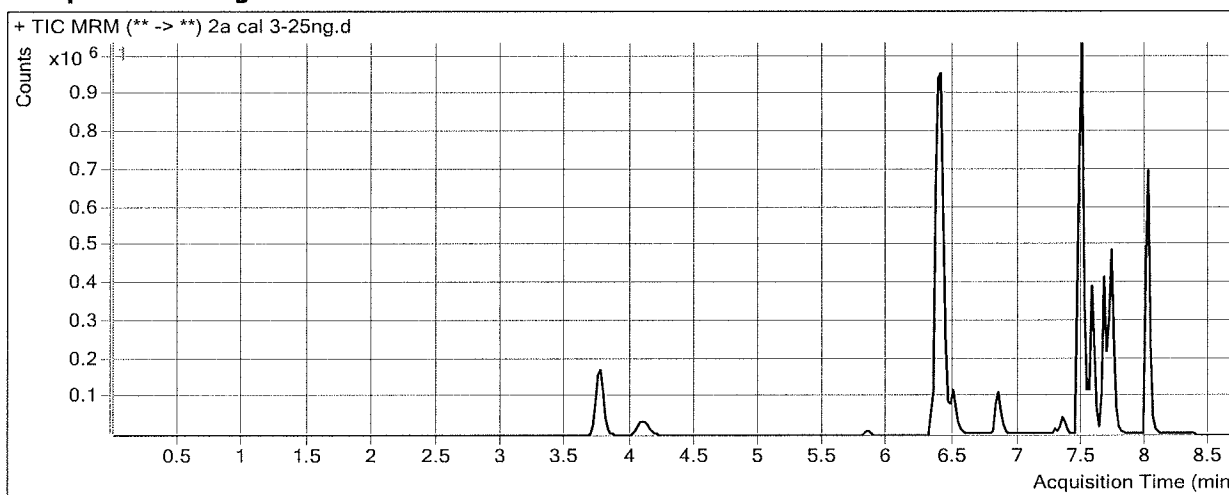
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 06:07 **Data File** 2a cal 3-25ng.d
Sample Type Calibration **Sample Name** 2a cal 3-25ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-C5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 252484 | 508831 | 0.4962 | 24.3825 |
| Levamisole | Cocaine-D3 | 4.084 | 138778 | 665680 | 0.2085 | 23.0103 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.851 | 18503 | 8226 | 2.2494 | 21.2778 |
| Zopiclone | Zopiclone-D4 | 6.388 | 64649 | 127403 | 0.5074 | 24.3920 |
| Cocaine | Cocaine-D3 | 6.352 | 390886 | 665680 | 0.5872 | 25.1127 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 582019 | 1087198 | 0.5353 | 24.6529 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 90304 | 1087198 | 0.0831 | 25.6948 |
| Ondansetron | Zopiclone-D4 | 6.496 | 201780 | 127403 | 1.5838 | 20.1951 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 215211 | 1087198 | 0.1980 | 22.3434 |
| Mitragynine | Doxepin-D3 | 7.272 | 10704 | 74404 | 0.1439 | 25.7175 |
| Doxepin | Doxepin-D3 | 7.344 | 34906 | 74404 | 0.4691 | 25.5796 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 852746 | 1837457 | 0.4641 | 25.0605 |
| Protriptyline | Protriptyline-D3 | 7.531 | 16203 | 35478 | 0.4567 | 22.9844 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 9806 | 26916 | 0.3643 | 23.9556 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 33820 | 26916 | 1.2565 | 24.5455 |
| Estazolam | Estazolam-D5 | 7.563 | 205997 | 398485 | 0.5170 | 23.5026 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 14124 | 26952 | 0.5240 | 22.3075 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 232846 | 577156 | 0.4034 | 25.7731 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.723 | 108443 | 468911 | 0.2313 | 27.3775 |
| Midazolam | Midazolam-D4 | 7.721 | 70977 | 411838 | 0.1723 | 25.0025 |
| Phenazepam | Phenazepam-D4 | 7.739 | 18012 | 26487 | 0.6800 | 25.5647 |
| Prazepam | Prazepam-D5 | 8.004 | 535562 | 1026927 | 0.5215 | 25.2494 |

ISP FORENSICS - Cd'A Instrument # 62340

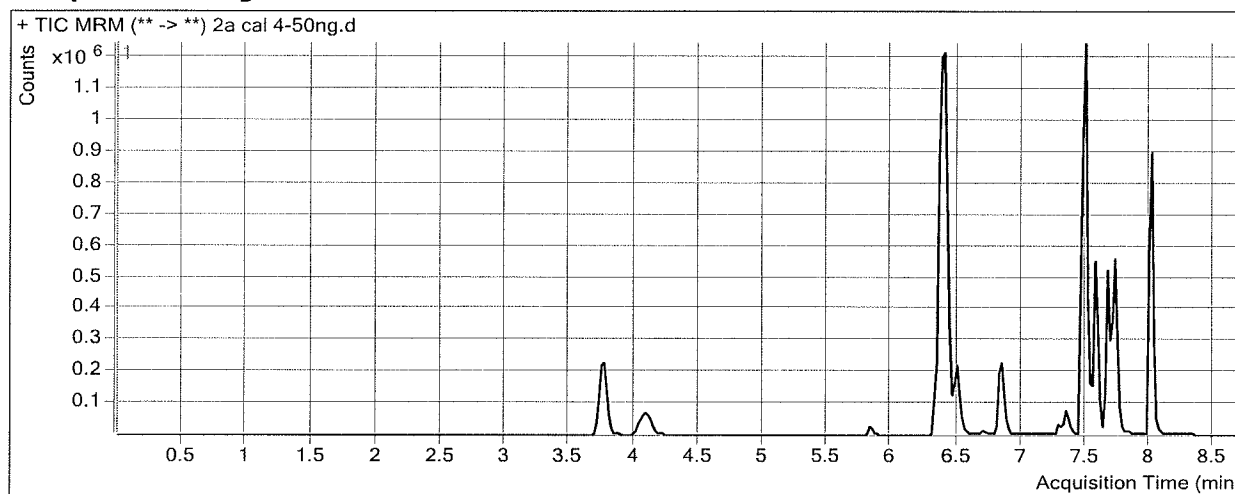
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 06:32 **Data File** 2a cal 4-50ng.d
Sample Type Calibration **Sample Name** 2a cal 4-50ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-D5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 498420 | 482637 | 1.0327 | 50.4235 |
| Levamisole | Cocaine-D3 | 4.084 | 271459 | 628126 | 0.4322 | 45.0230 |
| Benzoyllecgonine | Benzoyllecgonine-D8 | 5.851 | 37088 | 7618 | 4.8682 | 44.0368 |
| Zopiclone | Zopiclone-D4 | 6.388 | 108619 | 102755 | 1.0571 | 50.6647 |
| Cocaine | Cocaine-D3 | 6.352 | 753212 | 628126 | 1.1991 | 51.3776 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 1164284 | 1065397 | 1.0928 | 50.2394 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 180560 | 1065397 | 0.1695 | 53.1645 |
| Ondansetron | Zopiclone-D4 | 6.496 | 407575 | 102755 | 3.9665 | 46.0010 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 456340 | 1065397 | 0.4283 | 45.5233 |
| Mitragynine | Doxepin-D3 | 7.272 | 26064 | 87917 | 0.2965 | 51.5648 |
| Doxepin | Doxepin-D3 | 7.344 | 84135 | 87917 | 0.9570 | 51.8768 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 1544863 | 1675784 | 0.9219 | 49.9320 |
| Protriptyline | Protriptyline-D3 | 7.531 | 40253 | 40010 | 1.0061 | 48.9002 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 26546 | 33298 | 0.7972 | 54.1899 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 83930 | 33298 | 2.5206 | 50.5671 |
| Estazolam | Estazolam-D5 | 7.563 | 377852 | 338901 | 1.1149 | 48.7255 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 36228 | 32000 | 1.1321 | 47.1051 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 425447 | 514361 | 0.8271 | 52.7391 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.723 | 189665 | 419697 | 0.4519 | 55.5000 |
| Midazolam | Midazolam-D4 | 7.721 | 137429 | 398690 | 0.3447 | 49.2535 |
| Phenazepam | Phenazepam-D4 | 7.739 | 30095 | 19871 | 1.5145 | 58.3936 |
| Prazepam | Prazepam-D5 | 8.004 | 1002964 | 951667 | 1.0539 | 51.6901 |

ISP FORENSICS - Cd'A Instrument # 62340

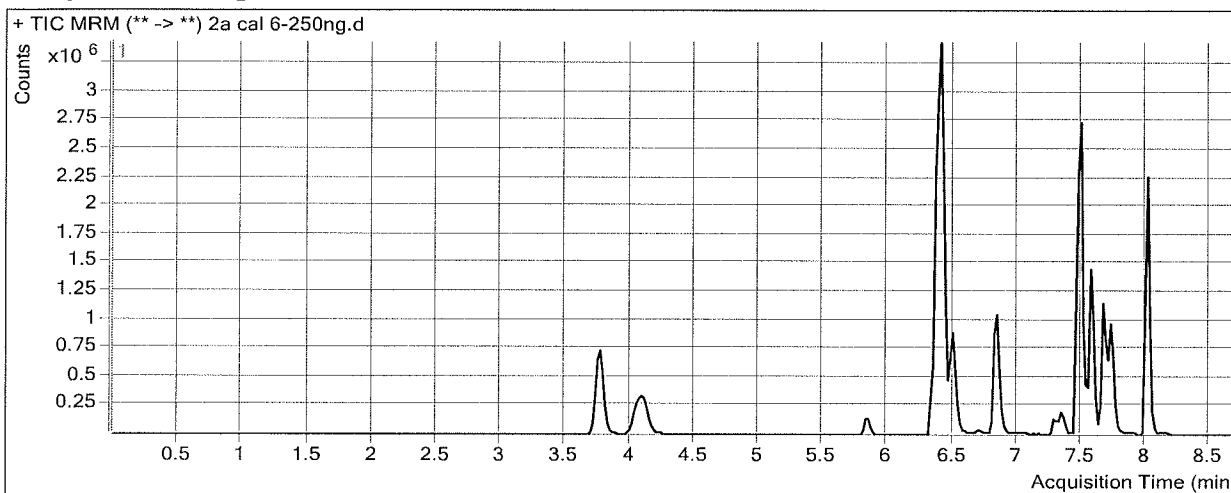
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 07:23 **Data File** 2a cal 6-250ng.d
Sample Type Calibration **Sample Name** 2a cal 6-250ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-F5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.770 | 2425085 | 483416 | 5.0166 | 243.7961 |
| Levamisole | Cocaine-D3 | 4.084 | 1341351 | 589238 | 2.2764 | 226.5039 |
| Benzoyllecgonine | Benzoyllecgonine-D8 | 5.851 | 215218 | 7828 | 27.4950 | 240.6787 |
| Zopiclone | Zopiclone-D4 | 6.388 | 407515 | 79226 | 5.1437 | 246.0104 |
| Cocaine | Cocaine-D3 | 6.352 | 3369078 | 589238 | 5.7177 | 245.3148 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 5077683 | 968996 | 5.2401 | 240.5892 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 734108 | 968996 | 0.7576 | 240.1138 |
| Ondansetron | Zopiclone-D4 | 6.496 | 1771842 | 79226 | 22.3643 | 245.2612 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 2109441 | 968996 | 2.1769 | 221.4616 |
| Mitragynine | Doxepin-D3 | 7.272 | 105544 | 71565 | 1.4748 | 251.1611 |
| Doxepin | Doxepin-D3 | 7.344 | 329933 | 71565 | 4.6102 | 248.8016 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 5913809 | 1284887 | 4.6026 | 249.9054 |
| Protriptyline | Protriptyline-D3 | 7.531 | 164858 | 32638 | 5.0511 | 239.7256 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 107532 | 26920 | 3.9945 | 277.4806 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 346618 | 26920 | 12.8758 | 263.7323 |
| Estazolam | Estazolam-D5 | 7.563 | 1345690 | 249755 | 5.3880 | 228.9643 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 150419 | 25384 | 5.9258 | 242.5922 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 1572611 | 407206 | 3.8620 | 245.8869 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 544553 | 272452 | 1.9987 | 252.6499 |
| Midazolam | Midazolam-D4 | 7.721 | 586108 | 332087 | 1.7649 | 249.0788 |
| Phenazepam | Phenazepam-D4 | 7.739 | 72513 | 11010 | 6.5859 | 257.9030 |
| Prazepam | Prazepam-D5 | 8.004 | 3807478 | 745756 | 5.1055 | 252.9130 |

ISP FORENSICS - Cd'A Instrument # 62340

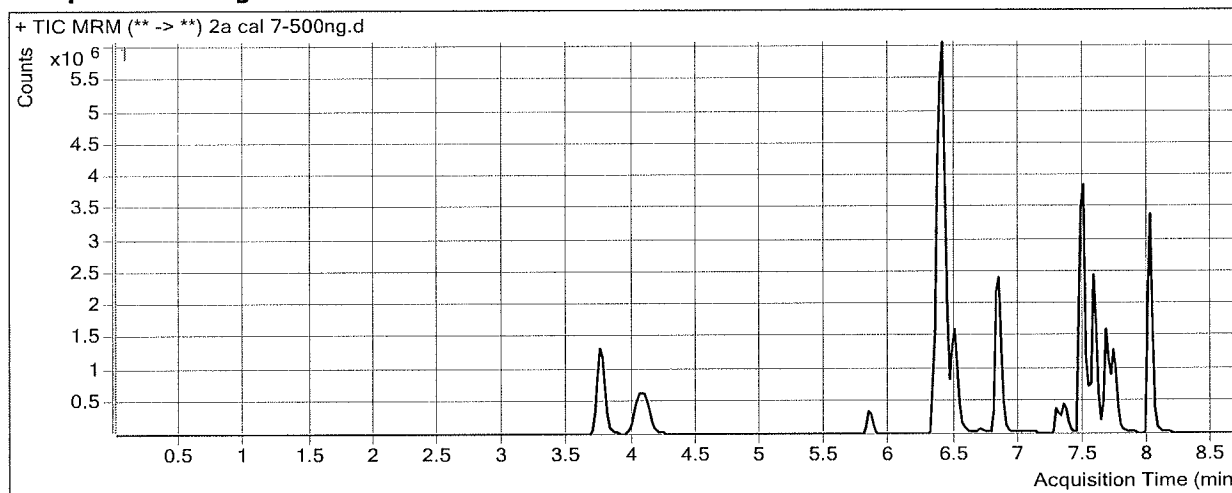
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 07:49 **Data File** 2a cal 7-500ng.d
Sample Type Calibration **Sample Name** 2a cal 7-500ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-G5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.750 | 4827063 | 464713 | 10.3872 | 504.4812 |
| Levamisole | Cocaine-D3 | 4.064 | 2734461 | 565181 | 4.8382 | 478.5943 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.851 | 524512 | 7436 | 70.5363 | 614.7358 |
| Zopiclone | Zopiclone-D4 | 6.388 | 555062 | 53125 | 10.4482 | 499.5729 |
| Cocaine | Cocaine-D3 | 6.352 | 6612012 | 565181 | 11.6989 | 502.0319 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 10205705 | 919306 | 11.1015 | 509.6085 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 1243594 | 919306 | 1.3528 | 429.2998 |
| Ondansetron | Zopiclone-D4 | 6.496 | 3411451 | 53125 | 64.2154 | 698.5346 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 5055893 | 919306 | 5.4997 | 555.7842 |
| Mitragynine | Doxepin-D3 | 7.272 | 302124 | 103414 | 2.9215 | 496.2174 |
| Doxepin | Doxepin-D3 | 7.344 | 908424 | 103414 | 8.7844 | 473.8058 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 9212581 | 1004758 | 9.1690 | 497.9959 |
| Protriptyline | Protriptyline-D3 | 7.531 | 368639 | 34412 | 10.7124 | 506.7936 |
| Maprotiline | Amitriptyline-D3 | 7.547 | 246734 | 36088 | 6.8370 | 475.9964 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 892652 | 36088 | 24.7353 | 507.8639 |
| Estazolam | Estazolam-D5 | 7.563 | 2136032 | 193304 | 11.0501 | 467.7907 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 352116 | 28011 | 12.5708 | 513.5814 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 2418611 | 323160 | 7.4843 | 476.4245 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 776426 | 203265 | 3.8198 | 484.7569 |
| Midazolam | Midazolam-D4 | 7.721 | 1021118 | 286437 | 3.5649 | 502.3339 |
| Phenazepam | Phenazepam-D4 | 7.739 | 93379 | 7990 | 11.6866 | 458.5671 |
| Prazepam | Prazepam-D5 | 8.004 | 6700960 | 663585 | 10.0981 | 500.8687 |

ISP FORENSICS - Cd'A Instrument # 62340

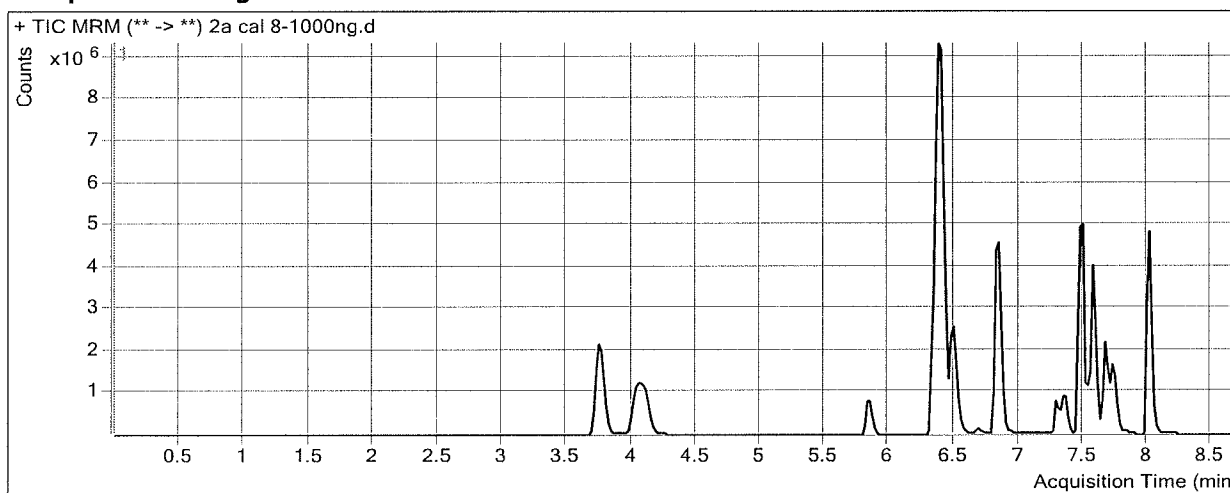
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2a.batch.bin
Analysis Time 12/21/2017 10:52 AM **Analyst Name** ISP Tox
Report Time 12/21/2017 10:55 AM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 10:52 AM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 08:14 **Data File** 2a cal 8-1000ng.d
Sample Type Calibration **Sample Name** 2a cal 8-1000ng
Dilution 1 **Acq Method** MDQ Panel 2a 3-12-17-3.m
Position P1-H5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.750 | 8826124 | 428082 | 20.6178 | 1001.0671 |
| Levamisole | Cocaine-D3 | 4.044 | 5429495 | 507665 | 10.6950 | 1054.9299 |
| Benzoylcegonine | Benzoylcegonine-D8 | 5.851 | 1338041 | 7971 | 167.8533 | 1460.4844 |
| Zopiclone | Zopiclone-D4 | 6.388 | 706329 | 33691 | 20.9650 | 1002.2896 |
| Cocaine | Cocaine-D3 | 6.352 | 11807836 | 507665 | 23.2591 | 998.1986 |
| Methylphenidate | Methylphenidate-D4 | 6.405 | 17481505 | 801722 | 21.8050 | 1000.8629 |
| Metoprolol | Methylphenidate-D4 | 6.421 | 1830915 | 801722 | 2.2837 | 725.2352 |
| Ondansetron | Zopiclone-D4 | 6.496 | 6084729 | 33691 | 180.6048 | 1959.1047 |
| Mirtazapine | Methylphenidate-D4 | 6.845 | 10167067 | 801722 | 12.6815 | 1278.3962 |
| Mitragynine | Doxepin-D3 | 7.272 | 624854 | 107210 | 5.8283 | 988.5941 |
| Doxepin | Doxepin-D3 | 7.364 | 2040887 | 107210 | 19.0363 | 1026.4276 |
| Carbamazepine | Carbamazepine-13C6 | 7.493 | 12286343 | 665485 | 18.4622 | 1002.8999 |
| Protriptyline | Protriptyline-D3 | 7.531 | 884389 | 41275 | 21.4266 | 1012.2286 |
| Maprotiline | Amitriptyline-D3 | 7.567 | 603787 | 50019 | 12.0711 | 841.5383 |
| Amitriptyline | Amitriptyline-D3 | 7.567 | 2253014 | 50019 | 45.0429 | 925.9022 |
| Estazolam | Estazolam-D5 | 7.563 | 2697825 | 107275 | 25.1487 | 1062.4672 |
| Nortriptyline | Nortriptyline-D3 | 7.571 | 894191 | 36728 | 24.3460 | 993.7805 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.659 | 3624961 | 260384 | 13.9216 | 886.1240 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.703 | 971159 | 65395 | 14.8507 | 1890.7238 |
| Midazolam | Midazolam-D4 | 7.721 | 1696944 | 238727 | 7.1083 | 1000.8907 |
| Phenazepam | Phenazepam-D4 | 7.739 | 129129 | 5764 | 22.4012 | 880.0865 |
| Prazepam | Prazepam-D5 | 8.004 | 10876077 | 545650 | 19.9323 | 989.2812 |

ISP FORENSICS - Cd'A Instrument # 62340

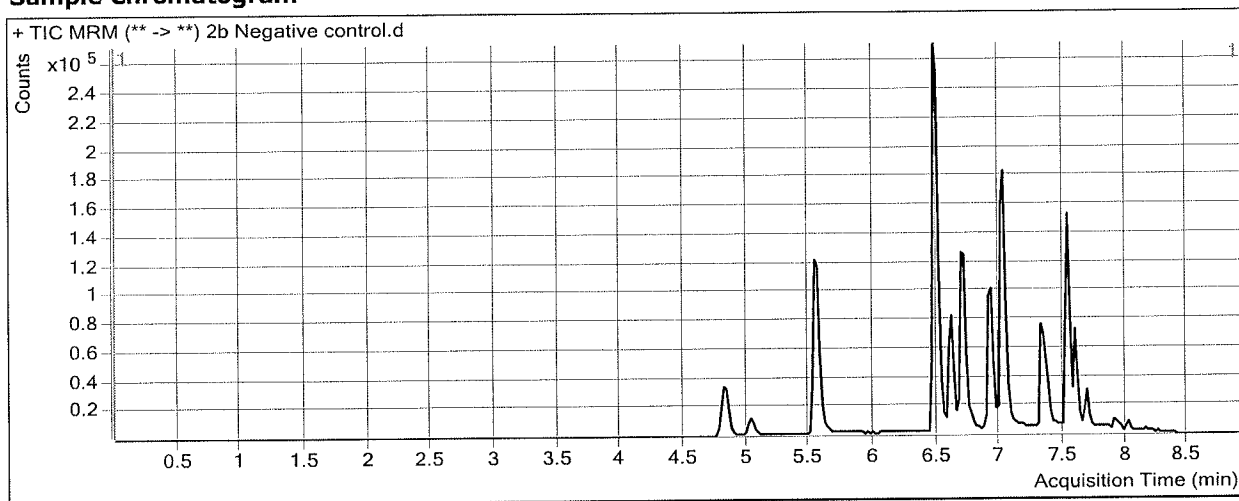
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 09:05 **Data File** 2b Negative control.d
Sample Type Sample **Sample Name** 2b Negative control
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-E6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



ISP FORENSICS - Cd'A Instrument # 62340

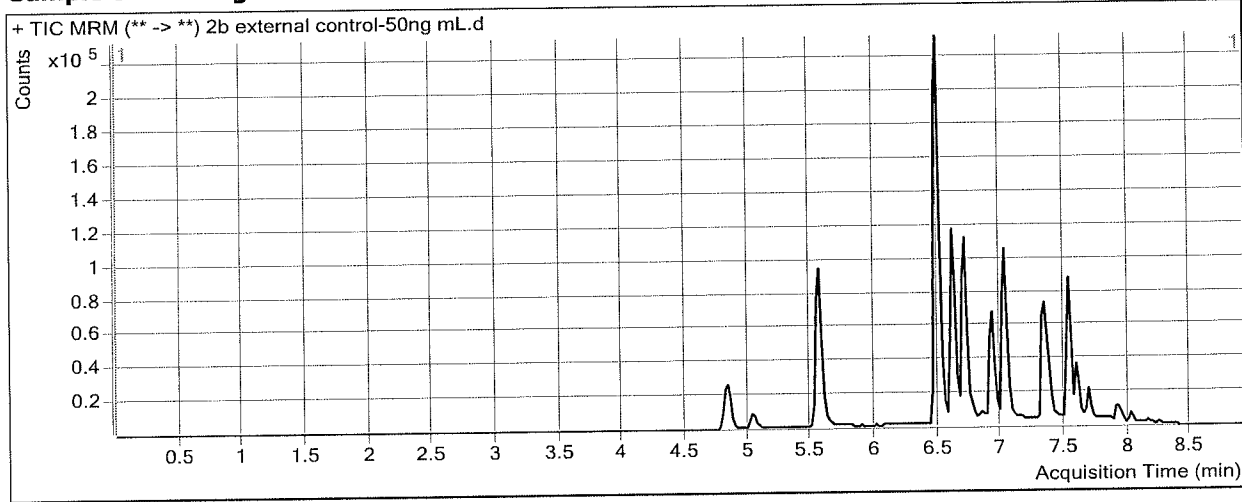
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 09:31 **Data File** 2b external control-50ng mL.d
Sample Type Sample **Sample Name** 2b external control-50ng mL
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-F6 **Sample Info**
Inj Vol -1 **Comment** AM 28 meperidine 50ng

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|------------|---------------|-------|----------|-----------|------------|------------|
| Meperidine | Meperidine-D4 | 6.616 | 120791 | 151349 | 0.7981 | 40.2656 |

ISP FORENSICS - Cd'A Instrument # 62340

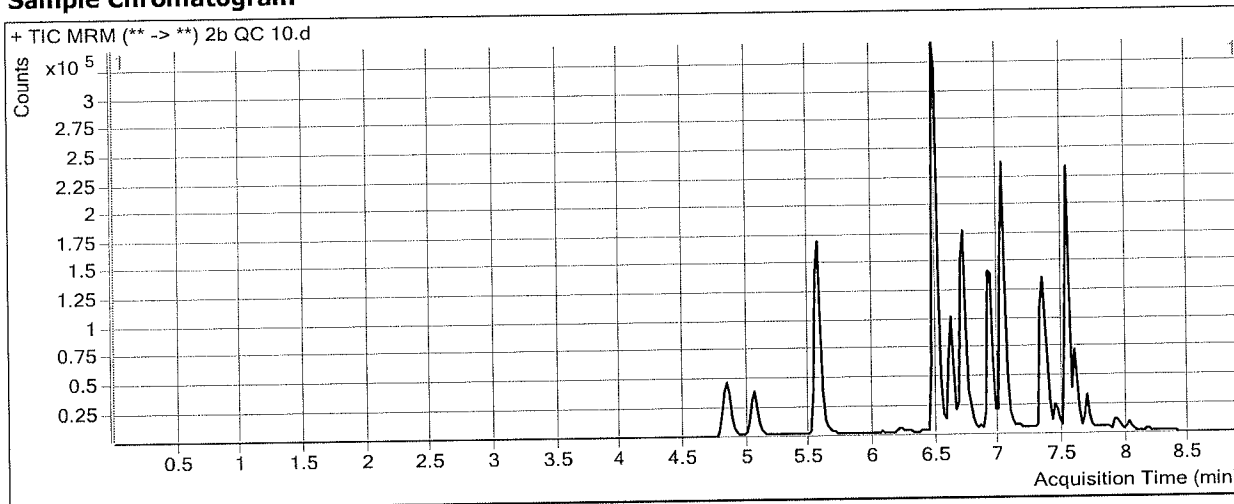
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 09:56 **Data File** 2b QC 10.d
Sample Type QC **Sample Name** 2b QC 10
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-A6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 49622 | 143211 | 0.3465 | 9.6269 |
| MDMA | MDMA-D6 | 5.048 | 66055 | 42063 | 1.5704 | 9.9555 |
| MDEA | MDEA-D6 | 5.548 | 75031 | 429255 | 0.1748 | 10.4734 |
| Primidone | Tapentadol-D3 | 6.215 | 7376 | 881343 | 0.0084 | 9.2640 |
| Tapentadol | Tapentadol-D3 | 6.483 | 188218 | 881343 | 0.2136 | 10.2233 |
| Meperidine | Meperidine-D4 | 6.616 | 44486 | 226747 | 0.1962 | 10.3134 |
| Normeperidine | Normeperidine-D4 | 6.724 | 49470 | 164763 | 0.3003 | 10.3728 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 52712 | 252795 | 0.2085 | 9.5589 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 86759 | 297533 | 0.2916 | 9.9612 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 103768 | 519875 | 0.1996 | 10.8067 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 69946 | 65041 | 1.0754 | 9.6194 |
| Zaleplon | Zaleplon-D4 | 7.341 | 53381 | 193755 | 0.2755 | 10.4566 |
| Phenytoin | Phenytoin-D10 | 7.389 | 2633 | 15673 | 0.1680 | 9.1033 |
| Verapamil | Trimipramine-D3 | 7.415 | 37148 | 107186 | 0.3466 | 10.7297 |
| Amoxapine | Clomipramine-D3 | 7.500 | 5456 | 46925 | 0.1163 | 10.2738 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 58731 | 74308 | 0.7904 | 9.5814 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 55705 | 65041 | 0.8565 | 10.0085 |
| Imipramine | Imipramine-D3 | 7.528 | 26975 | 83160 | 0.3244 | 10.2249 |
| Desipramine | Desipramine-D3 | 7.549 | 16043 | 83298 | 0.1926 | 9.8521 |
| Trimipramine | Trimipramine-D3 | 7.586 | 23804 | 107186 | 0.2221 | 10.8797 |
| Clomipramine | Clomipramine-D3 | 7.681 | 8817 | 46925 | 0.1879 | 9.8434 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 6064 | 19801 | 0.3062 | 11.6489 |

ISP FORENSICS - Cd'A Instrument # 62340

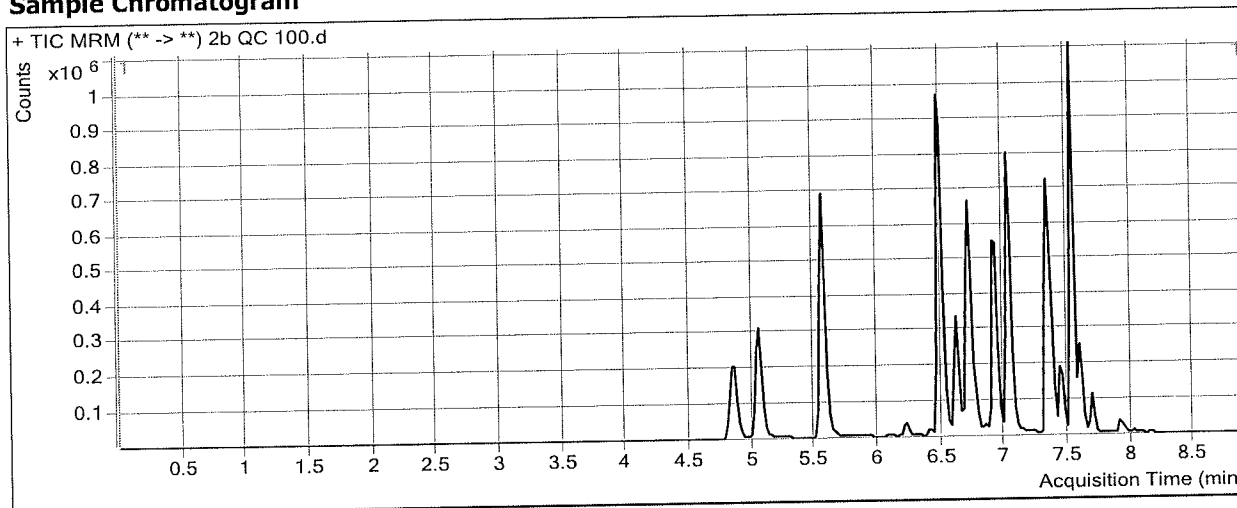
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 13:32 **Data File** 2b QC 100.d
Sample Type QC **Sample Name** 2b QC 100
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-B6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 492821 | 139551 | 3.5315 | 105.1172 |
| MDMA | MDMA-D6 | 5.048 | 676284 | 42536 | 15.8990 | 102.4712 |
| MDEA | MDEA-D6 | 5.548 | 774729 | 434985 | 1.7810 | 103.2446 |
| Primidone | Tapentadol-D3 | 6.215 | 75046 | 837375 | 0.0896 | 94.7294 |
| Tapentadol | Tapentadol-D3 | 6.483 | 1789083 | 837375 | 2.1365 | 101.4104 |
| Meperidine | Meperidine-D4 | 6.616 | 471833 | 231265 | 2.0402 | 102.0775 |
| Normeperidine | Normeperidine-D4 | 6.724 | 484131 | 167655 | 2.8877 | 104.5290 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 532597 | 253212 | 2.1034 | 98.6110 |
| Pentazocine | Pentazocine-13C3 | 6.899 | 860765 | 285547 | 3.0144 | 102.5575 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 1086049 | 525745 | 2.0657 | 107.3714 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 777844 | 50319 | 15.4583 | 93.2612 |
| Zaleplon | Zaleplon-D4 | 7.341 | 504093 | 176148 | 2.8618 | 101.6509 |
| Phenytoin | Phenytoin-D10 | 7.389 | 30288 | 15168 | 1.9968 | 99.0598 |
| Verapamil | Trimipramine-D3 | 7.415 | 372275 | 121173 | 3.0723 | 102.6333 |
| Amoxapine | Clomipramine-D3 | 7.500 | 49138 | 55122 | 0.8914 | 79.5308 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 636517 | 77441 | 8.2194 | 99.6537 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 420617 | 50319 | 8.3590 | 81.7043 |
| Imipramine | Imipramine-D3 | 7.528 | 308467 | 88920 | 3.4690 | 98.3636 |
| Desipramine | Desipramine-D3 | 7.549 | 176377 | 91198 | 1.9340 | 95.7716 |
| Trimipramine | Trimipramine-D3 | 7.586 | 253069 | 121173 | 2.0885 | 96.8756 |
| Clomipramine | Clomipramine-D3 | 7.681 | 93238 | 55122 | 1.6915 | 96.6970 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 63378 | 21234 | 2.9848 | 93.6871 |

ISP FORENSICS - Cd'A Instrument # 62340

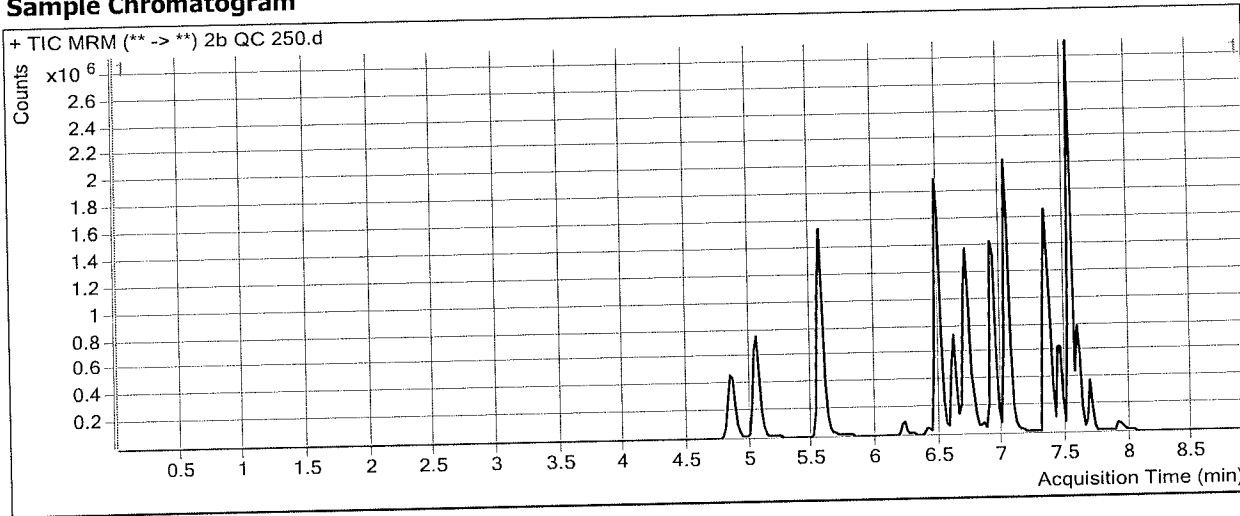
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 13:58 **Data File** 2b QC 250.d
Sample Type QC **Sample Name** 2b QC 250
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-C6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|-----------------------|-----------------------------|------------------|--------------------|------------------|--------------------|--|
| MDA | MDA-D5 | 4.850 | 1203923 | 140880 | 8.5458 | 255.4524 |
| MDMA | MDMA-D6 | 5.048 | 1720057 | 43343 | 39.6844 | 256.0464 |
| MDEA | MDEA-D6 | 5.548 | 2012604 | 445509 | 4.5175 | 261.2941 |
| Primidone | Tapentadol-D3 | 6.215 | 197187 | 773691 | 0.2549 | 268.5456 |
| Tapentadol | Tapentadol-D3 | 6.483 | 4139319 | 773691 | 5.3501 | 253.7962 |
| Meperidine | Meperidine-D4 | 6.616 | 1246469 | 251678 | 4.9526 | 247.0062 |
| Normeperidine | Normeperidine-D4 | 6.703 | 1245254 | 174117 | 7.1518 | 259.7025 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 1251863 | 234255 | 5.3440 | 250.9118 |
| Pentazocine | Pentazocine-13C3 | 6.899 | 2368593 | 335361 | 7.0628 | 240.2312 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 3164635 | 634321 | 4.9890 | 258.6395 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 2293207 | 33622 | 68.2046 | 400.0001 <i>outside curve range</i> |
| Zaleplon | Zaleplon-D4 | 7.341 | 1039599 | 144388 | 7.2000 | 254.6236 |
| Phenytoin | Phenytoin-D10 | 7.389 | 70456 | 15255 | 4.6186 | 228.0175 |
| Verapamil | Trimipramine-D3 | 7.415 | 1365426 | 183499 | 7.4411 | 249.9394 |
| Amoxapine | Clomipramine-D3 | 7.500 | 164072 | 85813 | 1.9120 | 170.7086 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 2089460 | 103141 | 20.2584 | 245.6183 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 792299 | 33622 | 23.5646 | 227.0106 |
| Imipramine | Imipramine-D3 | 7.528 | 1085208 | 111203 | 9.7588 | 274.6545 |
| Desipramine | Desipramine-D3 | 7.549 | 611046 | 120927 | 5.0530 | 249.6622 |
| Trimipramine | Trimipramine-D3 | 7.586 | 959735 | 183499 | 5.2302 | 241.6308 |
| Clomipramine | Clomipramine-D3 | 7.681 | 371026 | 85813 | 4.3236 | 248.7396 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 159738 | 19993 | 7.9898 | 246.9771 |

ISP FORENSICS - Cd'A Instrument # 62340

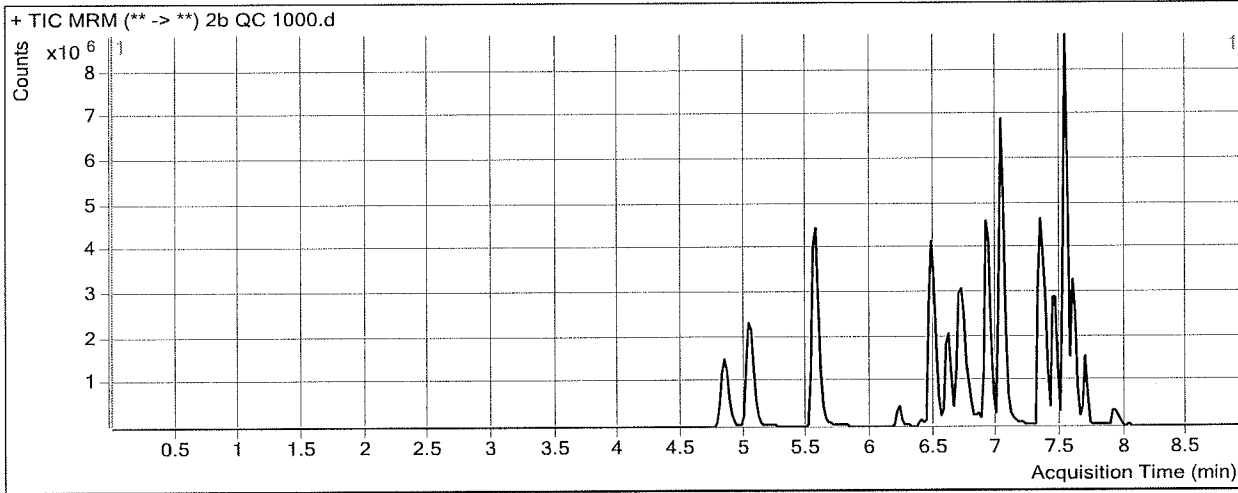
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 14:23 **Data File** 2b QC 1000.d
Sample Type QC **Sample Name** 2b QC 1000
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-D6 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.830 | 4159893 | 125037 | 33.2693 | 996.7000 |
| MDMA | MDMA-D6 | 5.028 | 6049041 | 38728 | 156.1936 | 1008.3103 |
| MDEA | MDEA-D6 | 5.548 | 6838147 | 390590 | 17.5072 | 1011.5289 |
| Primidone | Tapentadol-D3 | 6.215 | 888900 | 540274 | 1.6453 | 1731.0739 |
| Tapentadol | Tapentadol-D3 | 6.483 | 11457876 | 540274 | 21.2075 | 1005.7503 |
| Meperidine | Meperidine-D4 | 6.616 | 4298775 | 208079 | 20.6593 | 1028.6125 |
| Normeperidine | Normeperidine-D4 | 6.703 | 3574337 | 122661 | 29.1400 | 1059.8531 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 3989041 | 204992 | 19.4595 | 914.2941 |
| Pentazocine | Pentazocine-13C3 | 6.899 | 8478932 | 273823 | 30.9650 | 1053.0773 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 12204608 | 599941 | 20.3430 | 1053.1472 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 9080550 | 10312 | 880.5656 | 5124.1719 |
| Zaleplon | Zaleplon-D4 | 7.341 | 2426273 | 82561 | 29.3878 | 1036.9920 |
| Phenytoin | Phenytoin-D10 | 7.389 | 319147 | 14289 | 22.3352 | 1099.4536 |
| Verapamil | Trimipramine-D3 | 7.435 | 6503145 | 220341 | 29.5140 | 994.1863 |
| Amoxapine | Clomipramine-D3 | 7.500 | 474514 | 91866 | 5.1653 | 461.3705 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 8612836 | 109379 | 78.7432 | 954.7102 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 1162918 | 10312 | 112.7713 | 1079.4836 |
| Imipramine | Imipramine-D3 | 7.528 | 4525515 | 123533 | 36.6341 | 1027.9181 |
| Desipramine | Desipramine-D3 | 7.549 | 2112011 | 101629 | 20.7817 | 1025.7035 |
| Trimipramine | Trimipramine-D3 | 7.586 | 4599181 | 220341 | 20.8730 | 962.3788 |
| Clomipramine | Clomipramine-D3 | 7.681 | 1734117 | 91866 | 18.8765 | 1089.3671 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.908 | 906927 | 25668 | 35.3332 | 1084.4445 |

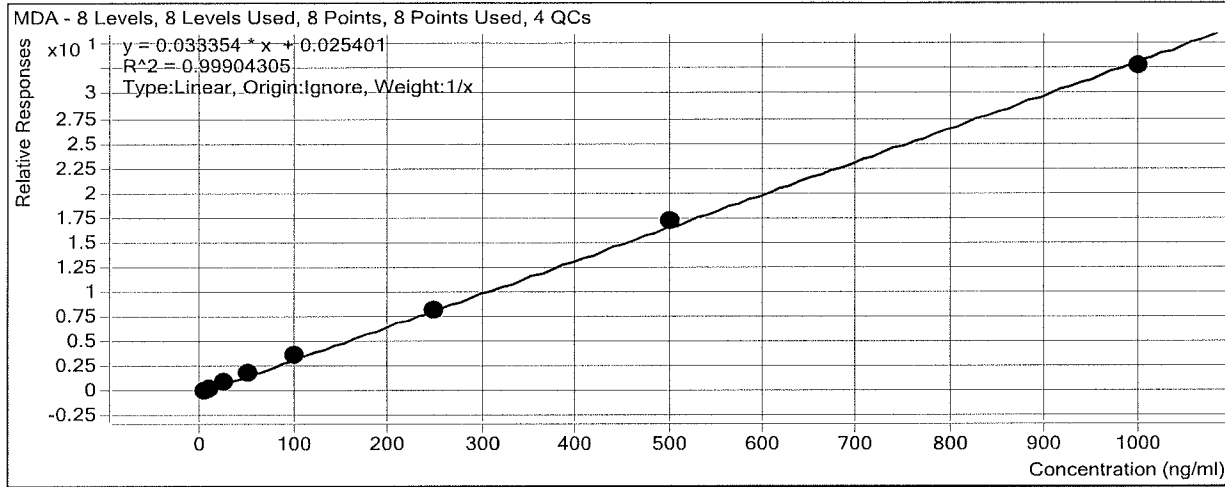
*outside
curve range*
*outside
curve
range*

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound MDA
Internal Standard MDA-D5



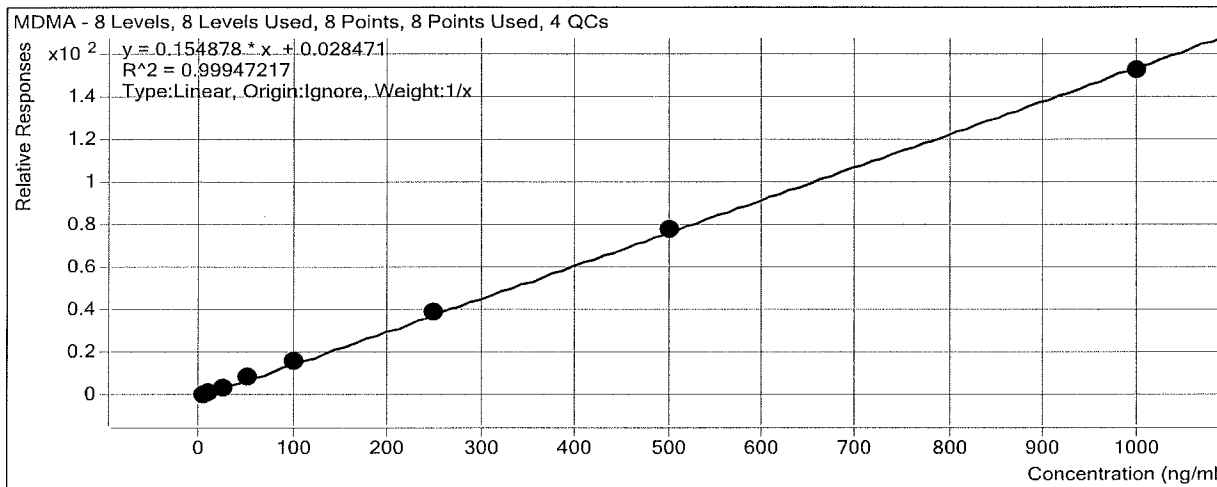
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.4 | 89.0 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.6 | 95.6 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.6 | 96.3 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.3 | 101.3 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 54.0 | 108.1 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 106.8 | 106.8 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 105.1 | 105.1 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 246.3 | 98.5 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 255.5 | 102.2 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 513.6 | 102.7 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 980.0 | 98.0 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 996.7 | 99.7 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound MDMA
Internal Standard MDMA-D6



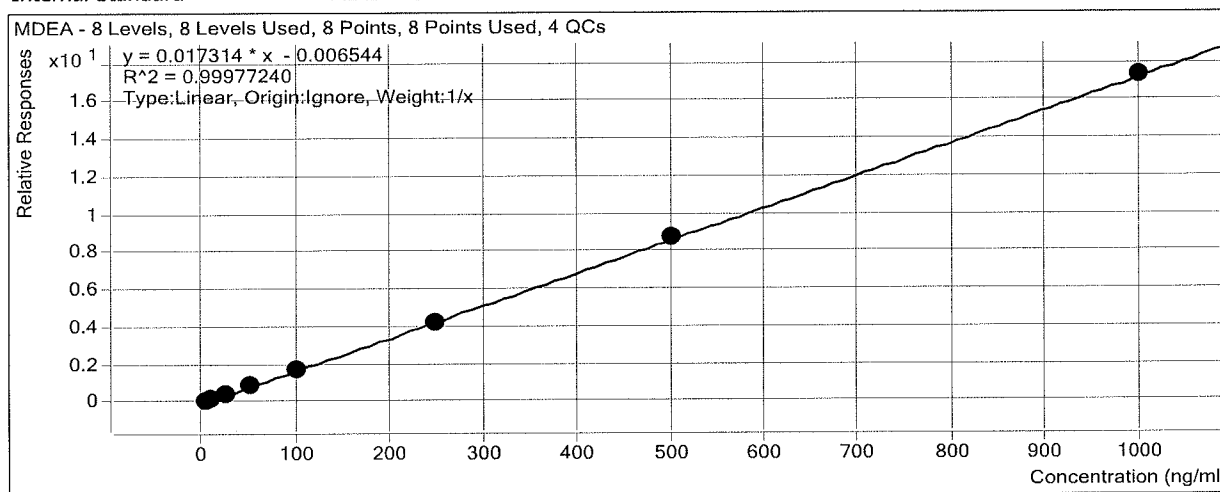
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.6 | 92.3 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.7 | 97.5 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 99.6 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.4 | 97.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 53.0 | 106.1 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 106.5 | 106.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 102.5 | 102.5 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 250.9 | 100.4 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 256.0 | 102.4 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 505.4 | 101.1 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 985.4 | 98.5 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1008.3 | 100.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound MDEA
Internal Standard MDEA-D6



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 103.9 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.6 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 104.7 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.8 | 95.2 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.9 | 103.8 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 100.1 | 100.1 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 103.2 | 103.2 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 242.8 | 97.1 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 261.3 | 104.5 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 505.3 | 101.1 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1001.0 | 100.1 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1011.5 | 101.2 |

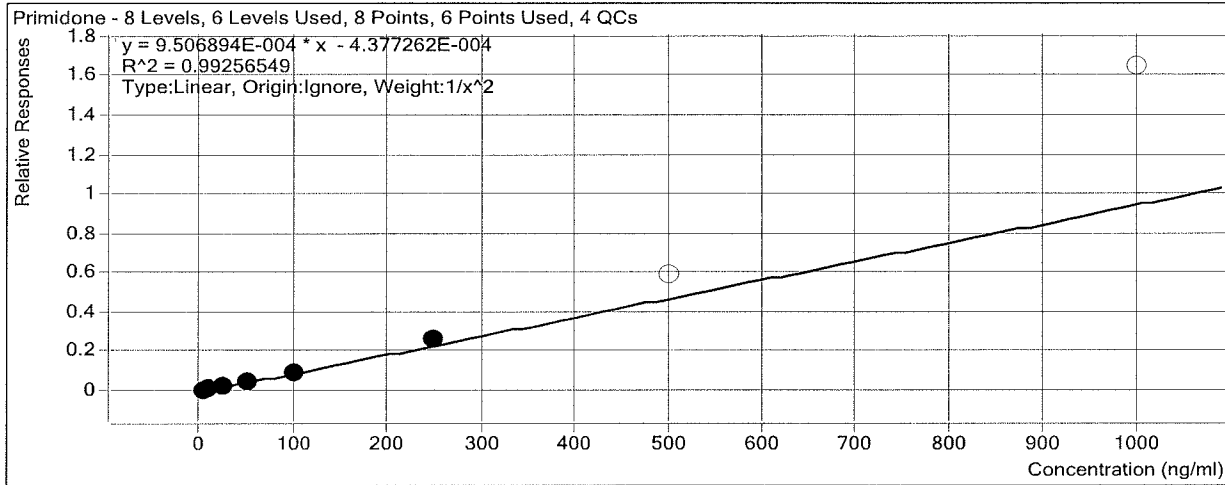
ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM

Analyst Name ISP TOX

Target Compound Primidone
Internal Standard Tapentadol-D3



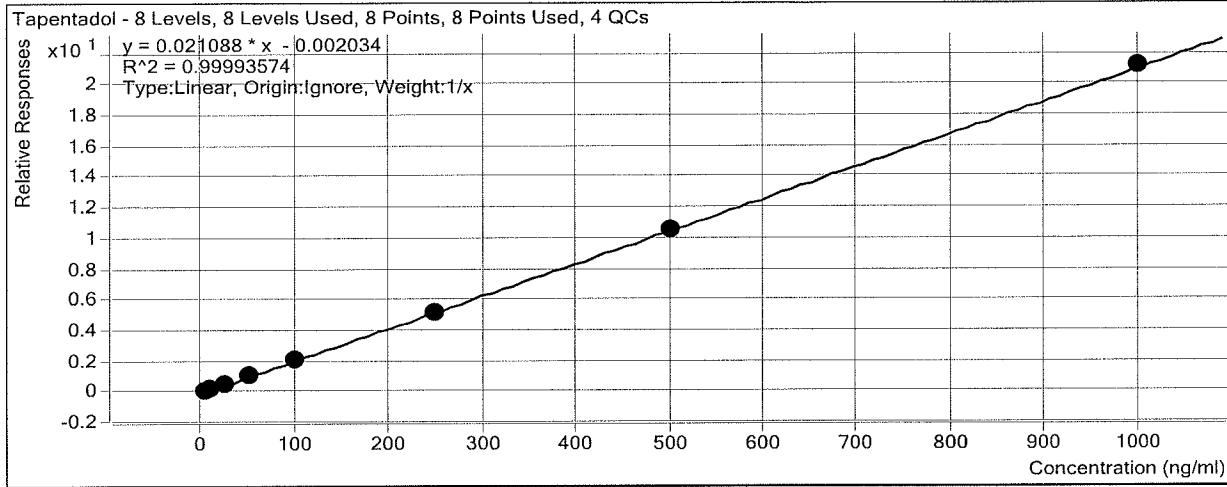
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 102.0 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 99.7 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.3 | 92.6 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.2 | 92.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 46.8 | 93.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 99.7 | 99.7 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 94.7 | 94.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 280.7 | 112.3 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 268.5 | 107.4 |
| 2b cal 7-500ng | 7 | <input type="checkbox"/> | 500 | 626.8 | 125.4 |
| 2b cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1730.9 | 173.1 |
| 2b QC 1000 | 8 | <input type="checkbox"/> | 1000 | 1731.1 | 173.1 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Tapentadol
Internal Standard Tapentadol-D3



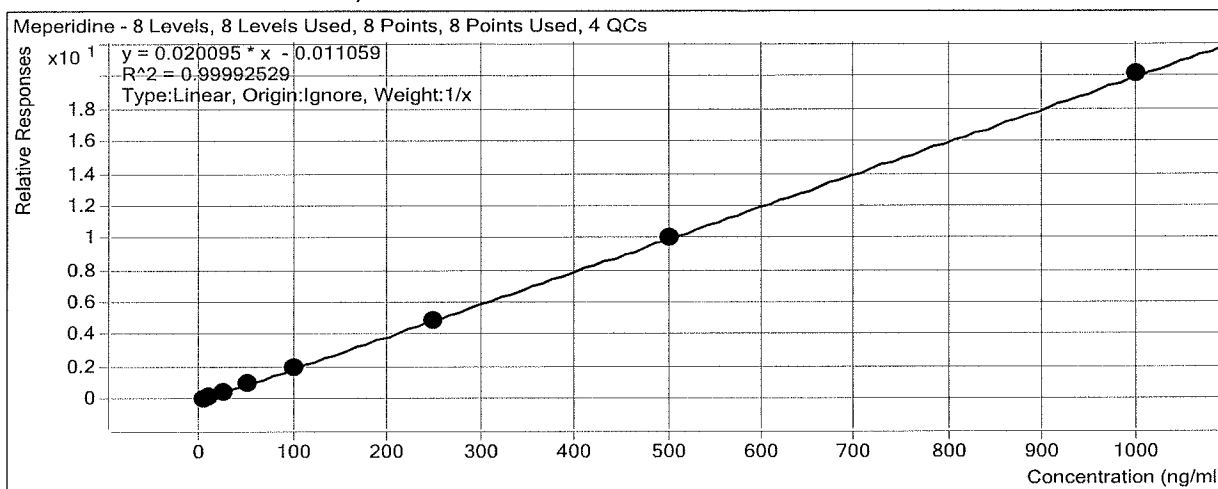
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 101.1 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.1 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.2 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.4 | 97.5 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.1 | 102.2 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 100.4 | 100.4 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 101.4 | 101.4 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 246.1 | 98.4 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 253.8 | 101.5 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 499.6 | 99.9 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1003.4 | 100.3 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1005.8 | 100.6 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Meperidine
Internal Standard Meperidine-D4



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 103.7 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.1 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.1 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.0 | 95.8 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 49.3 | 98.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 99.2 | 99.2 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 102.1 | 102.1 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 246.9 | 98.8 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 247.0 | 98.8 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 502.4 | 100.5 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1002.7 | 100.3 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1028.6 | 102.9 |

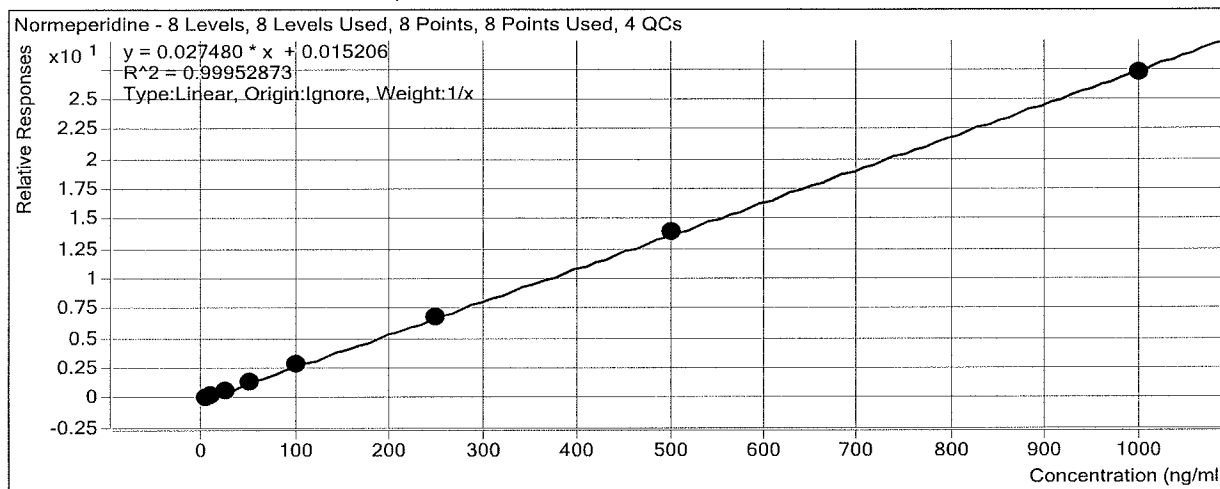
ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Normeperidine

Internal Standard Normeperidine-D4



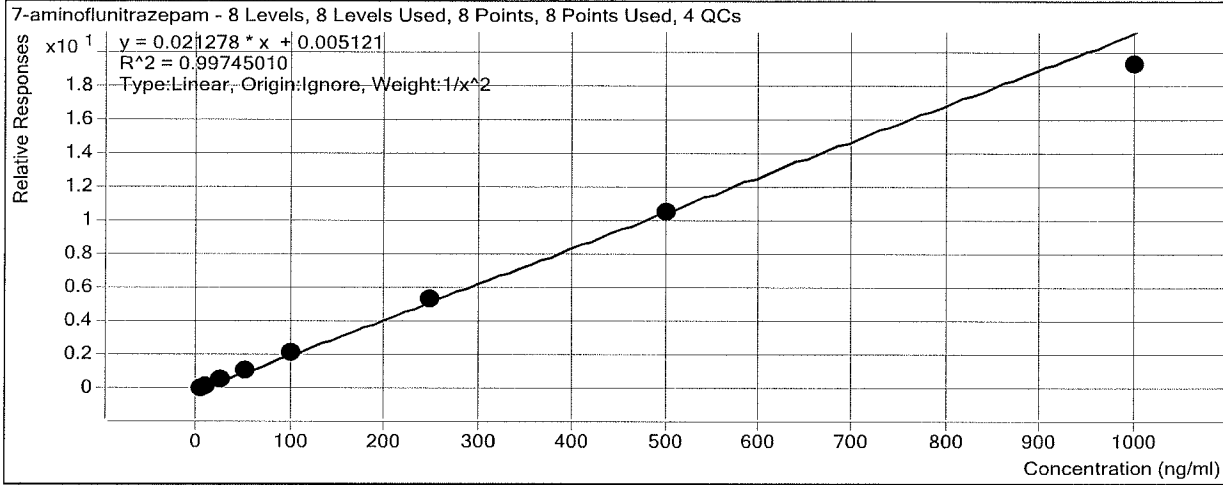
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.7 | 93.5 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.5 | 95.5 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.4 | 103.7 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.9 | 99.6 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 52.7 | 105.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 107.0 | 107.0 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 104.5 | 104.5 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 248.6 | 99.5 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 259.7 | 103.9 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 503.7 | 100.7 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 988.9 | 98.9 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1059.9 | 106.0 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound 7-aminoflunitrazepam
Internal Standard 7-aminoflunitrazepam-D7



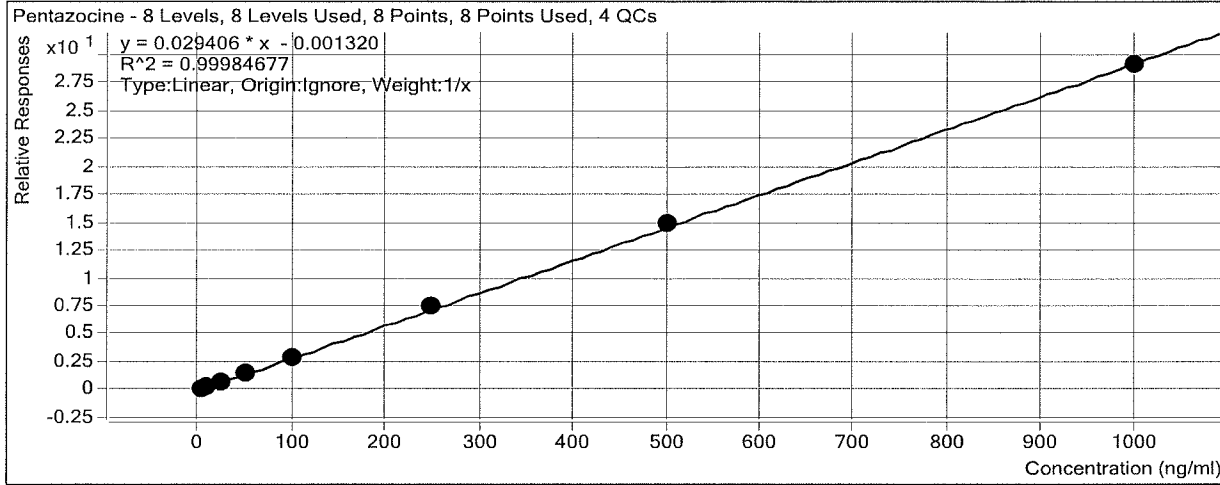
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.0 | 99.8 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.3 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.6 | 95.6 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.6 | 102.4 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 52.1 | 104.2 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 104.3 | 104.3 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 98.6 | 98.6 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 253.2 | 101.3 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 250.9 | 100.4 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 494.8 | 99.0 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 907.0 | 90.7 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 914.3 | 91.4 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Pentazocine
Internal Standard Pentazocine-13C3



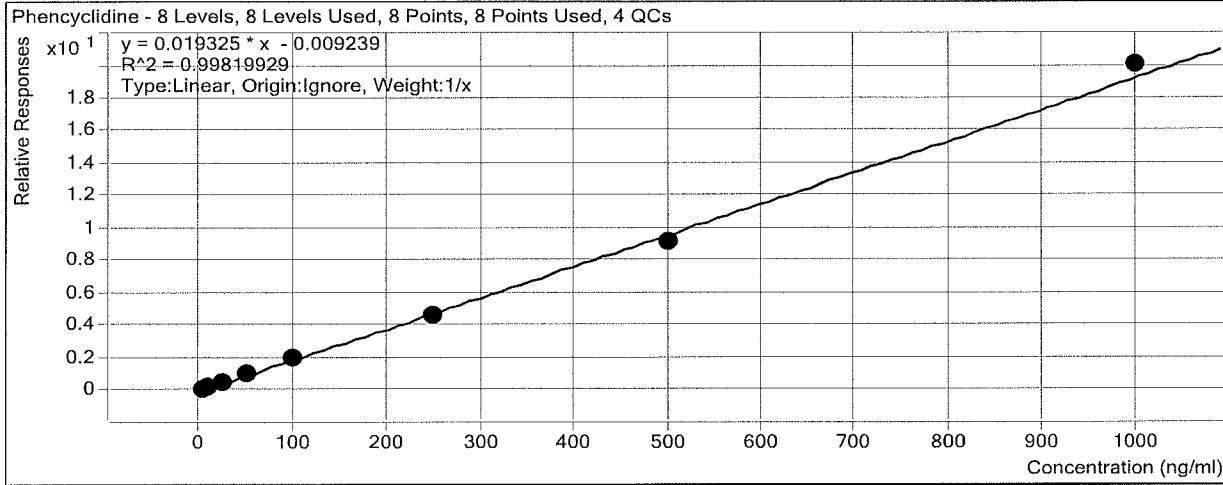
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.0 | 100.9 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 99.5 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 99.6 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.7 | 98.6 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 49.7 | 99.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 99.5 | 99.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 102.6 | 102.6 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 254.0 | 101.6 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 240.2 | 96.1 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 507.1 | 101.4 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 990.1 | 99.0 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1053.1 | 105.3 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Phencyclidine
Internal Standard Phencyclidine-D5



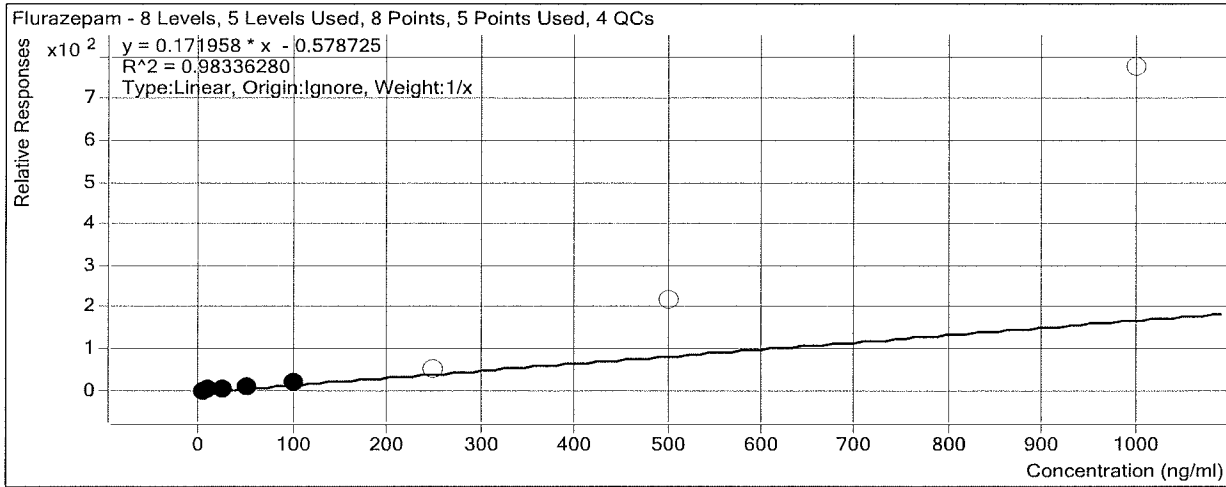
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 104.2 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 103.1 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.8 | 108.1 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.2 | 96.8 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.0 | 101.9 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 100.1 | 100.1 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 107.4 | 107.4 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 238.1 | 95.2 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 258.6 | 103.5 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 475.0 | 95.0 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1036.1 | 103.6 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1053.1 | 105.3 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Flurazepam*
Internal Standard *Flunitrazepam-D7*



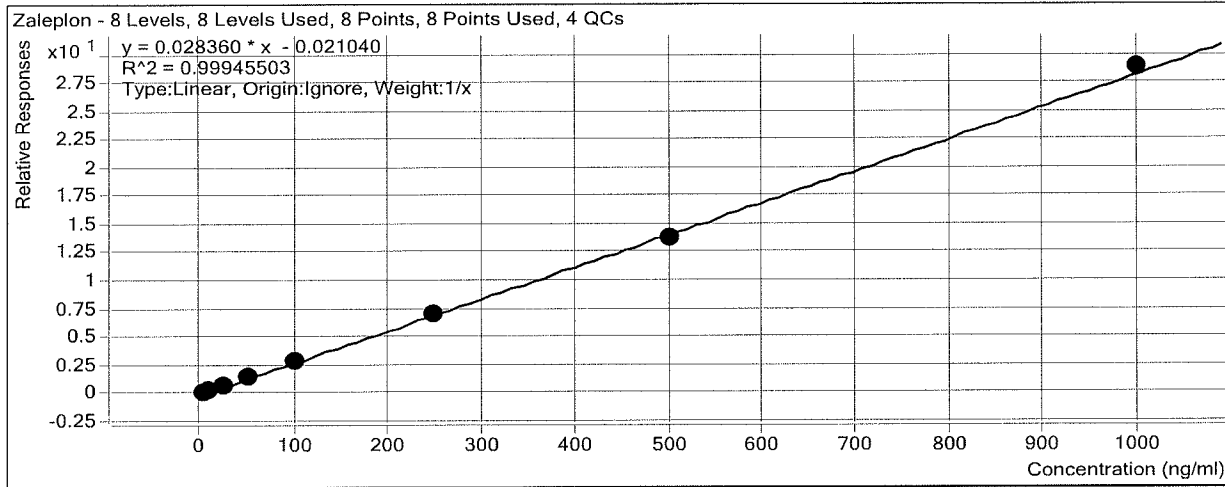
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.3 | 126.1 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.2 | 91.7 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.6 | 96.2 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 20.3 | 81.2 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 46.7 | 93.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 107.5 | 107.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 93.3 | 93.3 |
| 2b cal 6-250ng | 6 | <input type="checkbox"/> | 250 | 319.6 | 127.8 |
| 2b QC 250 | 6 | <input type="checkbox"/> | 250 | 400.0 | 160.0 |
| 2b cal 7-500ng | 7 | <input type="checkbox"/> | 500 | 1288.3 | 257.7 |
| 2b cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 4508.3 | 450.8 |
| 2b QC 1000 | 8 | <input type="checkbox"/> | 1000 | 5124.2 | 512.4 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Zaleplon*
Internal Standard *Zaleplon-D4*



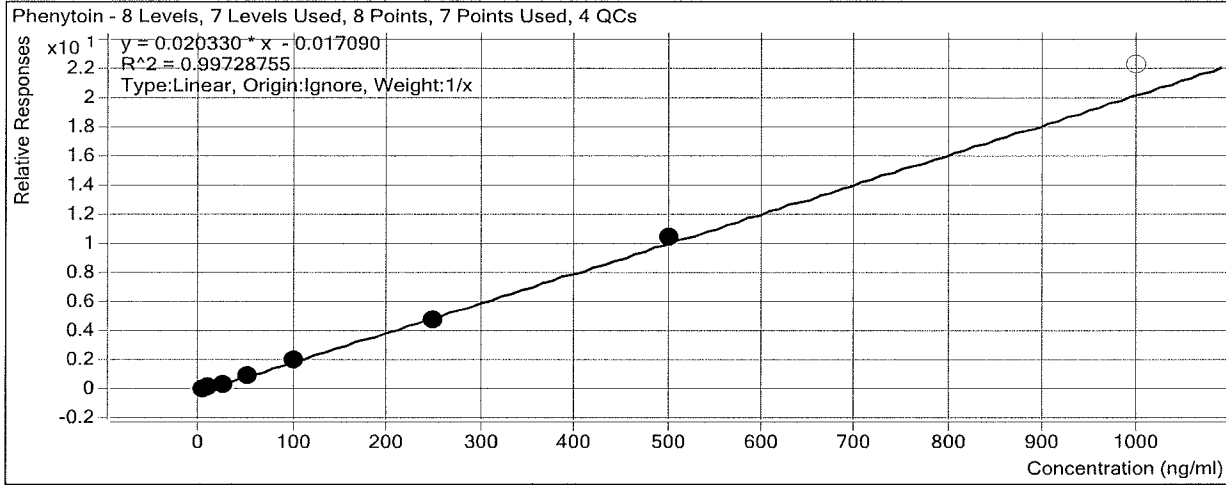
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.5 | 109.6 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.6 | 106.2 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 104.6 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 22.6 | 90.3 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 48.8 | 97.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 97.4 | 97.4 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 101.7 | 101.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 248.0 | 99.2 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 254.6 | 101.8 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 489.5 | 97.9 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1017.7 | 101.8 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1037.0 | 103.7 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Phenytoin
Internal Standard Phenytoin-D10



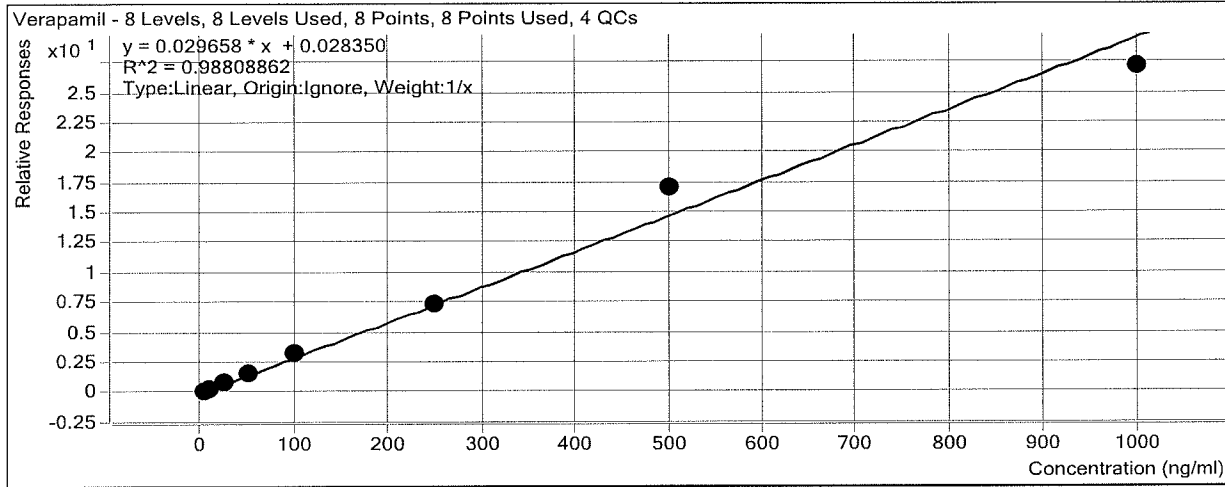
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.2 | 123.1 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.1 | 100.8 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.1 | 91.0 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 20.5 | 81.9 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 46.9 | 93.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 101.6 | 101.6 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 99.1 | 99.1 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 239.6 | 95.8 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 228.0 | 91.2 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 515.3 | 103.1 |
| 2b cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 1096.1 | 109.6 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1099.5 | 109.9 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Verapamil*
Internal Standard *Trimipramine-D3*



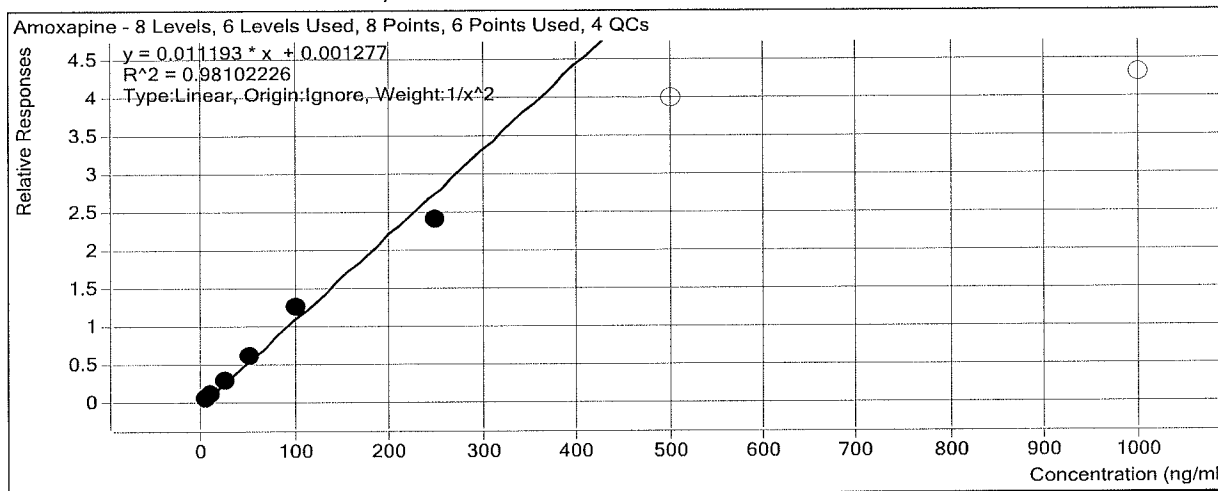
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.5 | 90.9 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 8.9 | 89.2 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.7 | 107.3 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.9 | 99.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.1 | 102.2 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 111.6 | 111.6 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 102.6 | 102.6 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 249.3 | 99.7 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 249.9 | 100.0 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 577.7 | 115.5 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 912.0 | 91.2 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 994.2 | 99.4 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound Amoxapine
Internal Standard Clomipramine-D3



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.3 | 105.5 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 8.5 | 85.4 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.3 | 102.7 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 26.1 | 104.4 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 53.0 | 105.9 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 112.5 | 112.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 79.5 | 79.5 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 215.8 | 86.3 |
| 2b QC 250 | 6 | <input type="checkbox"/> | 250 | 170.7 | 68.3 |
| 2b cal 7-500ng | 7 | <input type="checkbox"/> | 500 | 357.0 | 71.4 |
| 2b cal 8-1000ng | 8 | <input type="checkbox"/> | 1000 | 384.8 | 38.5 |
| 2b QC 1000 | 8 | <input type="checkbox"/> | 1000 | 461.4 | 46.1 |

ISP Forensics Calibration Curve Report

Batch Data Path

D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update

12/21/2017 12:27 PM

Analyst Name

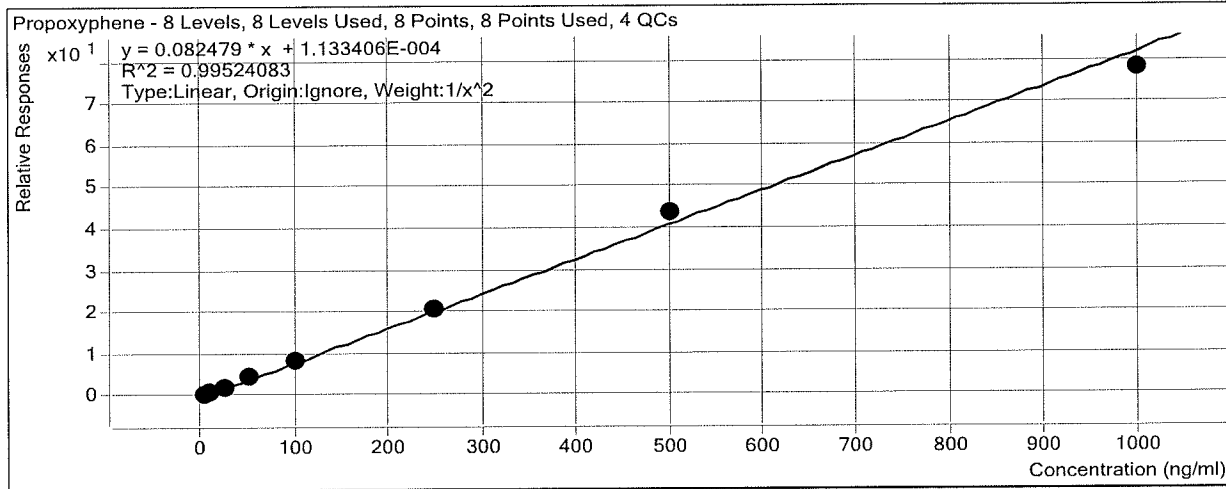
ISP TOX

Target Compound

Propoxyphene

Internal Standard

Propoxyphene-D11



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 104.3 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.0 | 89.8 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.6 | 95.8 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.2 | 100.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 54.0 | 107.9 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 98.5 | 98.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 99.7 | 99.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 246.9 | 98.8 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 245.6 | 98.2 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 528.2 | 105.6 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 944.6 | 94.5 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 954.7 | 95.5 |

ISP Forensics Calibration Curve Report

Batch Data Path

D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update

12/21/2017 12:27 PM

Analyst Name

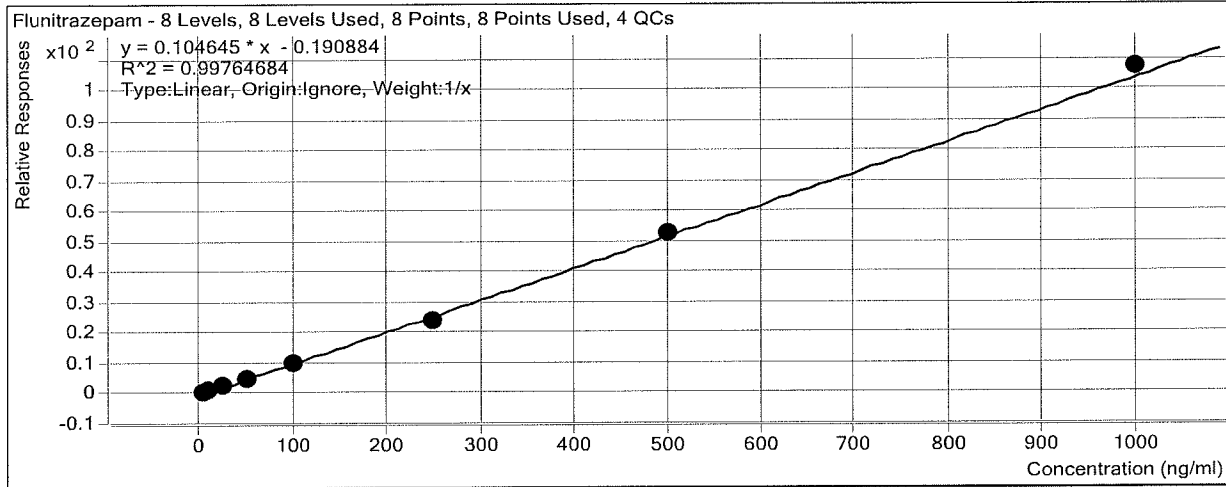
ISP TOX

Target Compound

Flunitrazepam

Internal Standard

Flunitrazepam-D7



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.3 | 125.2 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.4 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.0 | 100.1 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.2 | 96.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 43.2 | 86.5 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 93.5 | 93.5 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 81.7 | 81.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 229.5 | 91.8 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 227.0 | 90.8 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 506.0 | 101.2 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1027.1 | 102.7 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1079.5 | 107.9 |

ISP Forensics Calibration Curve Report

Batch Data Path

D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update

12/21/2017 12:27 PM

Analyst Name

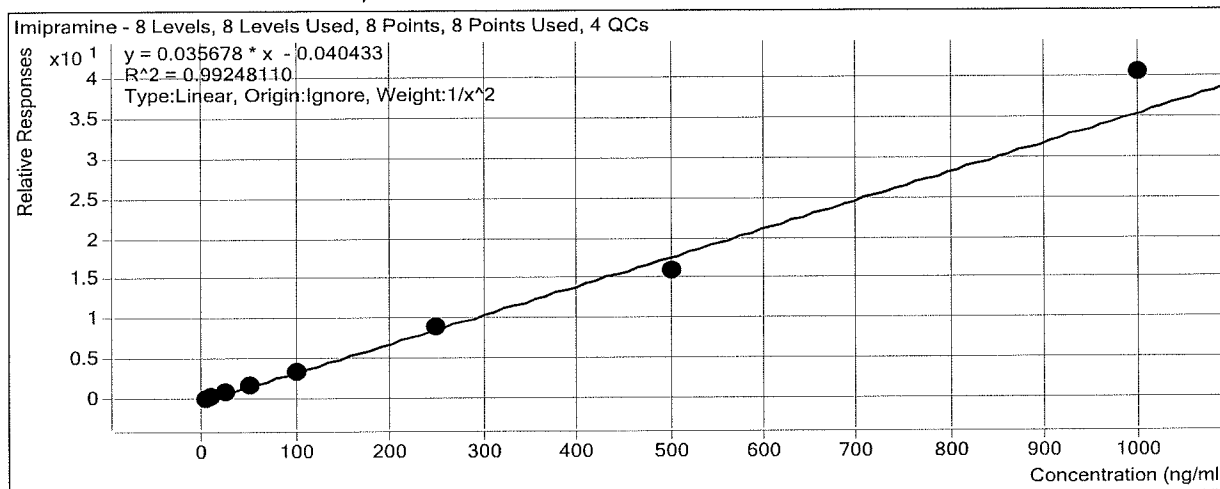
ISP TOX

Target Compound

Imipramine

Internal Standard

Imipramine-D3



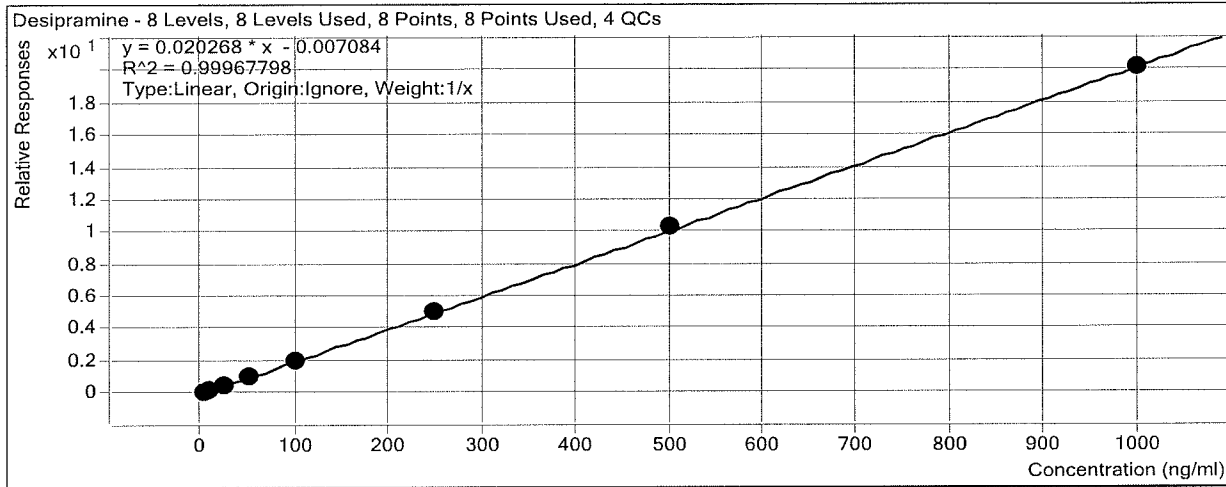
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.2 | 104.5 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.1 | 91.0 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.2 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.7 | 98.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 51.7 | 103.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 97.4 | 97.4 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 98.4 | 98.4 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 253.1 | 101.2 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 274.7 | 109.9 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 451.2 | 90.2 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1134.8 | 113.5 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1027.9 | 102.8 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Desipramine*
Internal Standard *Desipramine-D3*



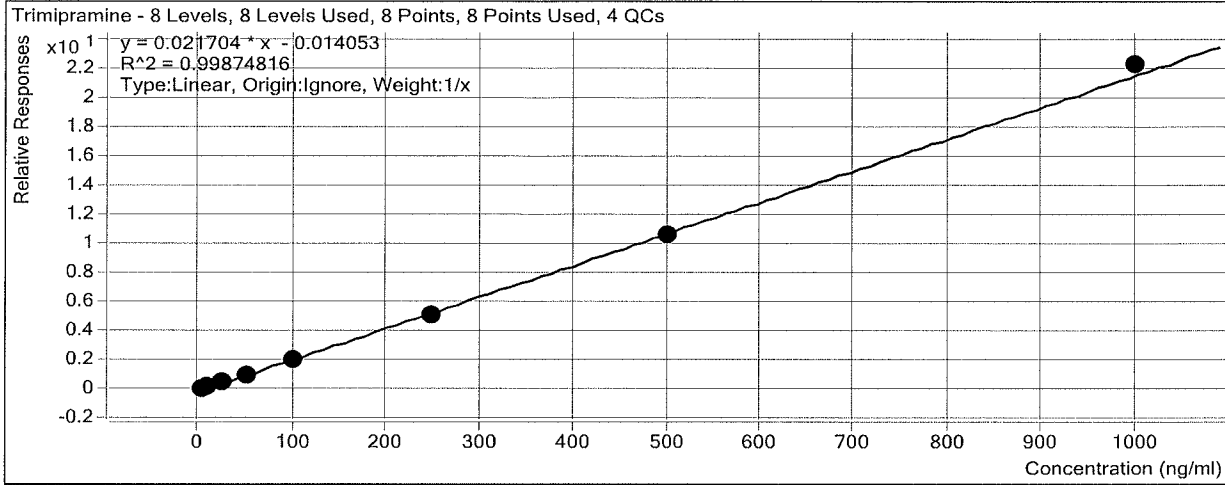
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 101.1 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.7 | 106.5 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.5 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 23.7 | 94.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 49.8 | 99.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 96.8 | 96.8 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 95.8 | 95.8 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 248.3 | 99.3 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 249.7 | 99.9 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 512.3 | 102.5 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 993.4 | 99.3 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1025.7 | 102.6 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Trimipramine*
Internal Standard *Trimipramine-D3*



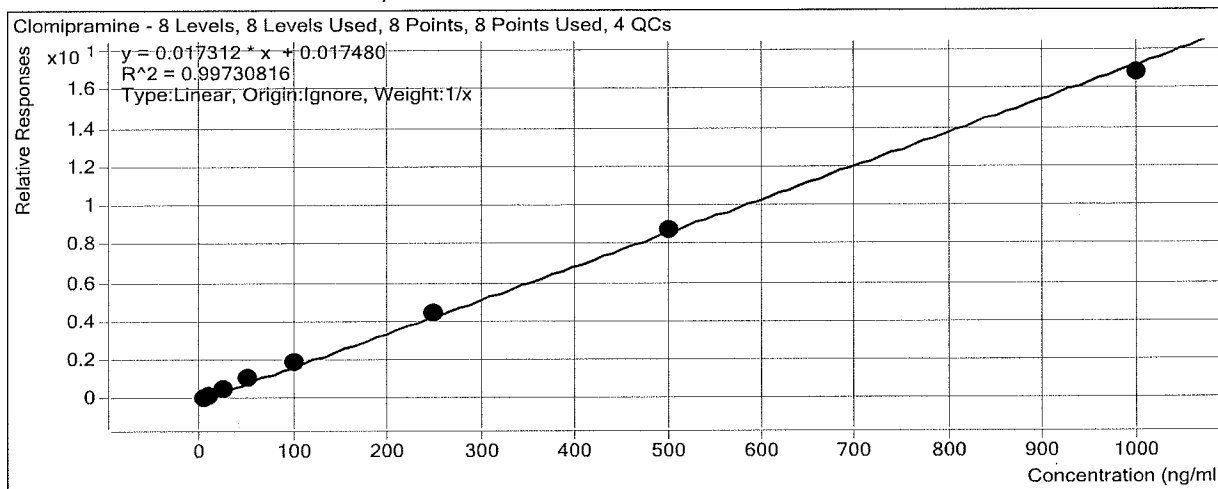
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.1 | 122.0 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 9.4 | 94.3 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 10.9 | 108.8 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 24.2 | 96.7 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 47.4 | 94.7 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 96.7 | 96.7 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 96.9 | 96.9 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 235.9 | 94.3 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 241.6 | 96.7 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 492.8 | 98.6 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1027.6 | 102.8 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 962.4 | 96.2 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Clomipramine*
Internal Standard *Clomipramine-D3*



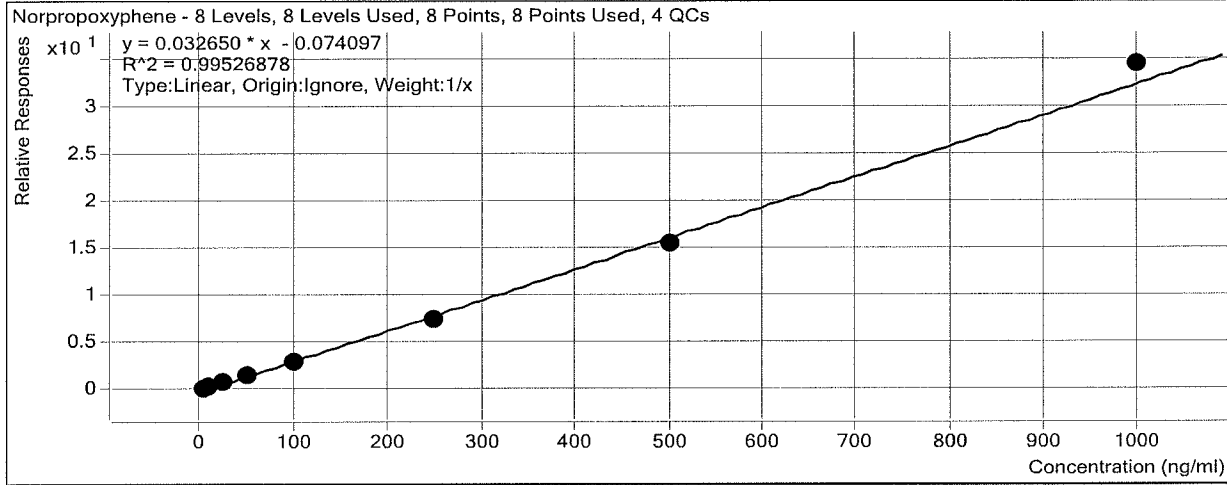
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 4.2 | 83.5 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 8.3 | 83.0 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.4 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 25.5 | 102.0 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 59.7 | 119.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 111.7 | 111.7 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 96.7 | 96.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 256.0 | 102.4 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 248.7 | 99.5 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 504.7 | 100.9 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 969.9 | 97.0 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1089.4 | 108.9 |

ISP Forensics Calibration Curve Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin

Last Calib Update 12/21/2017 12:27 PM **Analyst Name** ISP TOX

Target Compound *Norpropoxyphene*
Internal Standard *Norpropoxyphene-D5*



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------|-------|-------------------------------------|----------|------------|----------|
| 2b cal 1-5ng | 1 | <input checked="" type="checkbox"/> | 5 | 6.7 | 133.6 |
| 2b cal 2-10ng | 2 | <input checked="" type="checkbox"/> | 10 | 10.5 | 104.6 |
| 2b QC 10 | 2 | <input checked="" type="checkbox"/> | 10 | 11.6 | 116.5 |
| 2b cal 3-25ng | 3 | <input checked="" type="checkbox"/> | 25 | 22.0 | 87.9 |
| 2b cal 4-50ng | 4 | <input checked="" type="checkbox"/> | 50 | 44.2 | 88.4 |
| 2b cal 5-100ng | 5 | <input checked="" type="checkbox"/> | 100 | 93.2 | 93.2 |
| 2b QC 100 | 5 | <input checked="" type="checkbox"/> | 100 | 93.7 | 93.7 |
| 2b cal 6-250ng | 6 | <input checked="" type="checkbox"/> | 250 | 227.6 | 91.1 |
| 2b QC 250 | 6 | <input checked="" type="checkbox"/> | 250 | 247.0 | 98.8 |
| 2b cal 7-500ng | 7 | <input checked="" type="checkbox"/> | 500 | 478.0 | 95.6 |
| 2b cal 8-1000ng | 8 | <input checked="" type="checkbox"/> | 1000 | 1057.9 | 105.8 |
| 2b QC 1000 | 8 | <input checked="" type="checkbox"/> | 1000 | 1084.4 | 108.4 |

ISP FORENSICS - Cd'A Instrument # 62340

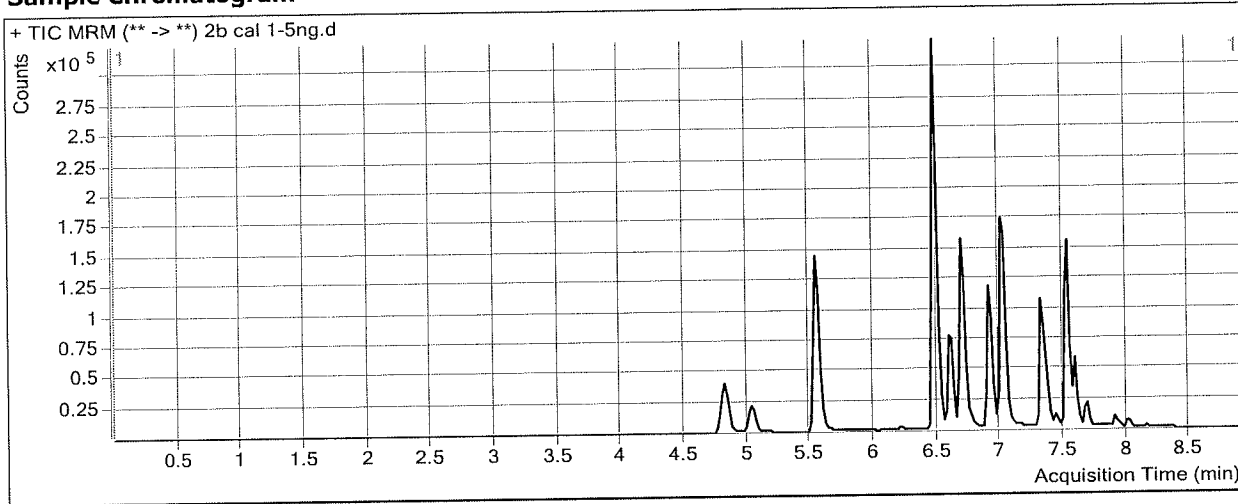
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 05:29 **Data File** 2b cal 1-5ng.d
Sample Type Calibration **Sample Name** 2b cal 1-5ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-A5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.830 | 24818 | 142820 | 0.1738 | 4.4482 |
| MDMA | MDMA-D6 | 5.048 | 30147 | 40557 | 0.7433 | 4.6156 |
| MDEA | MDEA-D6 | 5.548 | 34896 | 418448 | 0.0834 | 5.1945 |
| Primidone | Tapentadol-D3 | 6.195 | 3800 | 861114 | 0.0044 | 5.1021 |
| Tapentadol | Tapentadol-D3 | 6.483 | 90045 | 861114 | 0.1046 | 5.0550 |
| Meperidine | Meperidine-D4 | 6.616 | 20589 | 221104 | 0.0931 | 5.1842 |
| Normeperidine | Normeperidine-D4 | 6.703 | 23101 | 160782 | 0.1437 | 4.6753 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.716 | 27723 | 249194 | 0.1113 | 4.9878 |
| Pentazocine | Pentazocine-13C3 | 6.899 | 39932 | 271579 | 0.1470 | 5.0452 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 44915 | 491038 | 0.0915 | 5.2112 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 33129 | 65517 | 0.5056 | 6.3060 |
| Zaleplon | Zaleplon-D4 | 7.321 | 25692 | 191172 | 0.1344 | 5.4808 |
| Phenytoin | Phenytoin-D10 | 7.389 | 1598 | 14792 | 0.1080 | 6.1553 |
| Verapamil | Trimipramine-D3 | 7.415 | 15055 | 92289 | 0.1631 | 4.5446 |
| Amoxapine | Clomipramine-D3 | 7.500 | 2532 | 41979 | 0.0603 | 5.2741 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 29194 | 67885 | 0.4300 | 5.2127 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 30416 | 65517 | 0.4642 | 6.2604 |
| Imipramine | Imipramine-D3 | 7.528 | 11119 | 76123 | 0.1461 | 5.2273 |
| Desipramine | Desipramine-D3 | 7.529 | 7121 | 74635 | 0.0954 | 5.0571 |
| Trimipramine | Trimipramine-D3 | 7.586 | 10923 | 92289 | 0.1184 | 6.1006 |
| Clomipramine | Clomipramine-D3 | 7.681 | 3768 | 41979 | 0.0898 | 4.1753 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 2266 | 15739 | 0.1439 | 6.6781 |

ISP FORENSICS - Cd'A Instrument # 62340

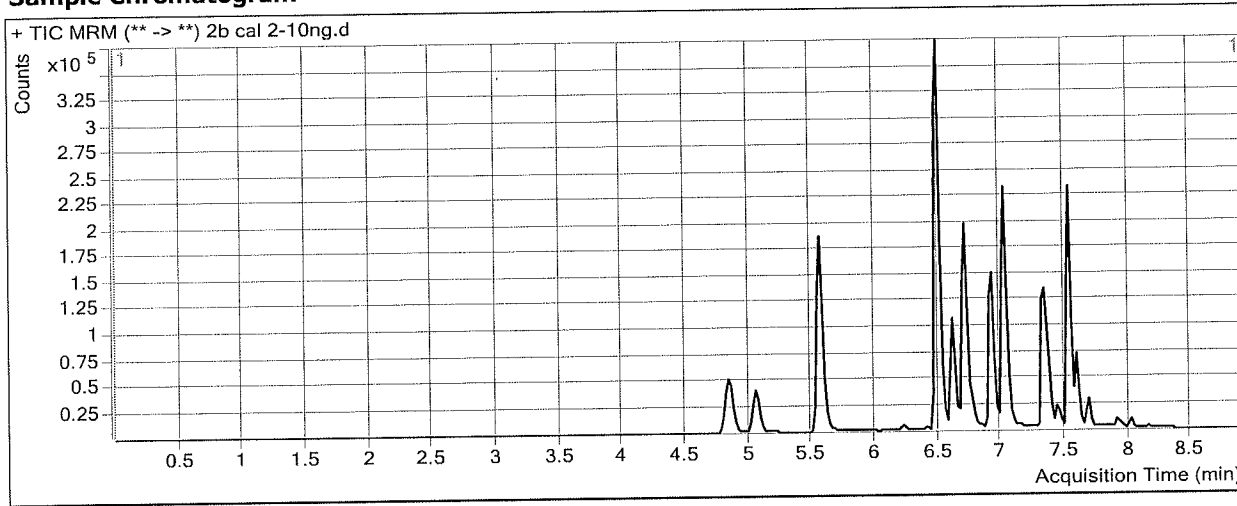
Multi-Drug Confirmatory Analysis Report

| | | | |
|--------------------------|---|----------------------|-----------|
| Batch Data Path | D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin | | |
| Analysis Time | 12/21/2017 12:27 PM | Analyst Name | ISP Tox |
| Report Time | 12/21/2017 12:27 PM | Reporter Name | ISP Tox |
| Last Calib Update | 12/21/2017 12:27 PM | Batch State | Processed |

Analysis Info

| | | | |
|--------------------|------------------|--------------------|-----------------------|
| Acq Time | 2017-12-20 05:54 | Data File | 2b cal 2-10ng.d |
| Sample Type | Calibration | Sample Name | 2b cal 2-10ng |
| Dilution | 1 | Acq Method | MDQ Panel 2b 8-8-17.m |
| Position | P1-B5 | Sample Info | |
| Inj Vol | -1 | Comment | AM 28 Panel 2 |

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 53938 | 156613 | 0.3444 | 9.5641 |
| MDMA | MDMA-D6 | 5.048 | 70132 | 45604 | 1.5378 | 9.7454 |
| MDEA | MDEA-D6 | 5.548 | 77299 | 470654 | 0.1642 | 9.8636 |
| Primidone | Tapentadol-D3 | 6.215 | 8572 | 948670 | 0.0090 | 9.9653 |
| Tapentadol | Tapentadol-D3 | 6.503 | 198385 | 948670 | 0.2091 | 10.0128 |
| Meperidine | Meperidine-D4 | 6.616 | 47534 | 242342 | 0.1961 | 10.3111 |
| Normeperidine | Normeperidine-D4 | 6.724 | 50626 | 182373 | 0.2776 | 9.5484 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 58583 | 273407 | 0.2143 | 9.8294 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 89025 | 305561 | 0.2913 | 9.9529 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 101059 | 531878 | 0.1900 | 10.3100 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 72848 | 72937 | 0.9988 | 9.1737 |
| Zaleplon | Zaleplon-D4 | 7.341 | 58928 | 210455 | 0.2800 | 10.6152 |
| Phenytoin | Phenytoin-D10 | 7.389 | 3016 | 16066 | 0.1878 | 10.0760 |
| Verapamil | Trimipramine-D3 | 7.415 | 32700 | 111672 | 0.2928 | 8.9173 |
| Amoxapine | Clomipramine-D3 | 7.500 | 4487 | 46326 | 0.0969 | 8.5405 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 57322 | 77404 | 0.7406 | 8.9773 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 64251 | 72937 | 0.8809 | 10.2422 |
| Imipramine | Imipramine-D3 | 7.528 | 24170 | 84997 | 0.2844 | 9.1033 |
| Desipramine | Desipramine-D3 | 7.529 | 17033 | 81551 | 0.2089 | 10.6548 |
| Trimipramine | Trimipramine-D3 | 7.586 | 21284 | 111672 | 0.1906 | 9.4292 |
| Clomipramine | Clomipramine-D3 | 7.681 | 7470 | 46326 | 0.1612 | 8.3044 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 4661 | 17433 | 0.2673 | 10.4573 |

ISP FORENSICS - Cd'A Instrument # 62340

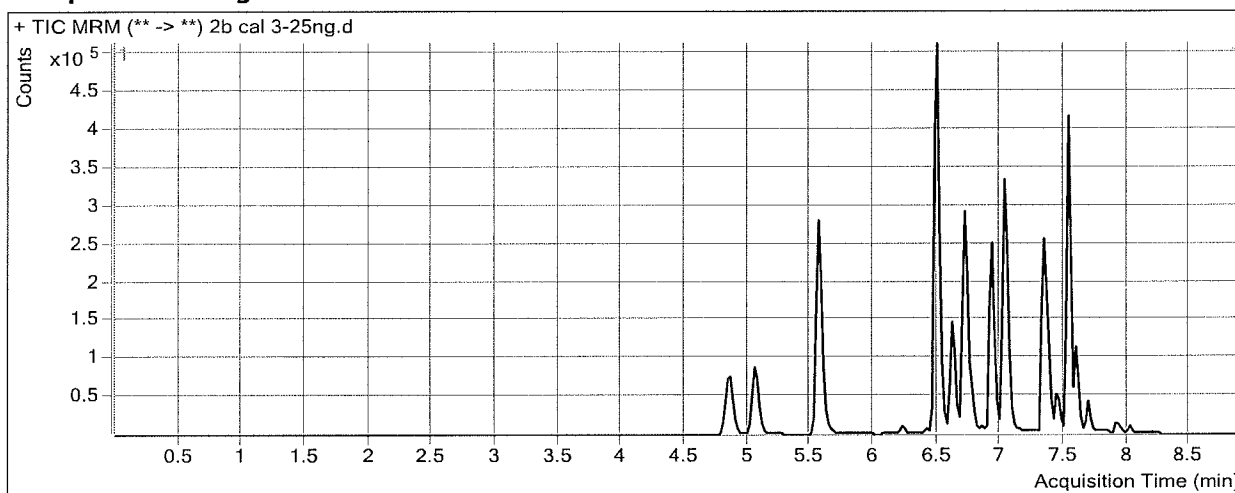
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 06:20 **Data File** 2b cal 3-25ng.d
Sample Type Calibration **Sample Name** 2b cal 3-25ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-C5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 137342 | 157840 | 0.8701 | 25.3264 |
| MDMA | MDMA-D6 | 5.048 | 176457 | 46314 | 3.8100 | 24.4162 |
| MDEA | MDEA-D6 | 5.568 | 195566 | 482023 | 0.4057 | 23.8108 |
| Primidone | Tapentadol-D3 | 6.215 | 20748 | 961207 | 0.0216 | 23.1659 |
| Tapentadol | Tapentadol-D3 | 6.503 | 492104 | 961207 | 0.5120 | 24.3737 |
| Meperidine | Meperidine-D4 | 6.616 | 120518 | 256198 | 0.4704 | 23.9591 |
| Normeperidine | Normeperidine-D4 | 6.724 | 128822 | 184235 | 0.6992 | 24.8915 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 151219 | 274919 | 0.5500 | 25.6099 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 230879 | 319044 | 0.7237 | 24.6545 |
| Phencyclidine | Phencyclidine-D5 | 7.048 | 265845 | 580098 | 0.4583 | 24.1920 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 191703 | 65817 | 2.9127 | 20.3038 |
| Zaleplon | Zaleplon-D4 | 7.341 | 133959 | 216293 | 0.6193 | 22.5806 |
| Phenytoin | Phenytoin-D10 | 7.389 | 7177 | 17970 | 0.3994 | 20.4869 |
| Verapamil | Trimipramine-D3 | 7.415 | 96129 | 125221 | 0.7677 | 24.9282 |
| Amoxapine | Clomipramine-D3 | 7.500 | 13053 | 44507 | 0.2933 | 26.0892 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 168302 | 81046 | 2.0766 | 25.1765 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 153958 | 65817 | 2.3392 | 24.1777 |
| Imipramine | Imipramine-D3 | 7.528 | 71560 | 85228 | 0.8396 | 24.6667 |
| Desipramine | Desipramine-D3 | 7.549 | 44075 | 93231 | 0.4728 | 23.6749 |
| Trimipramine | Trimipramine-D3 | 7.586 | 63914 | 125221 | 0.5104 | 24.1648 |
| Clomipramine | Clomipramine-D3 | 7.681 | 20429 | 44507 | 0.4590 | 25.5042 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 12722 | 19780 | 0.6432 | 21.9689 |

ISP FORENSICS - Cd'A Instrument # 62340

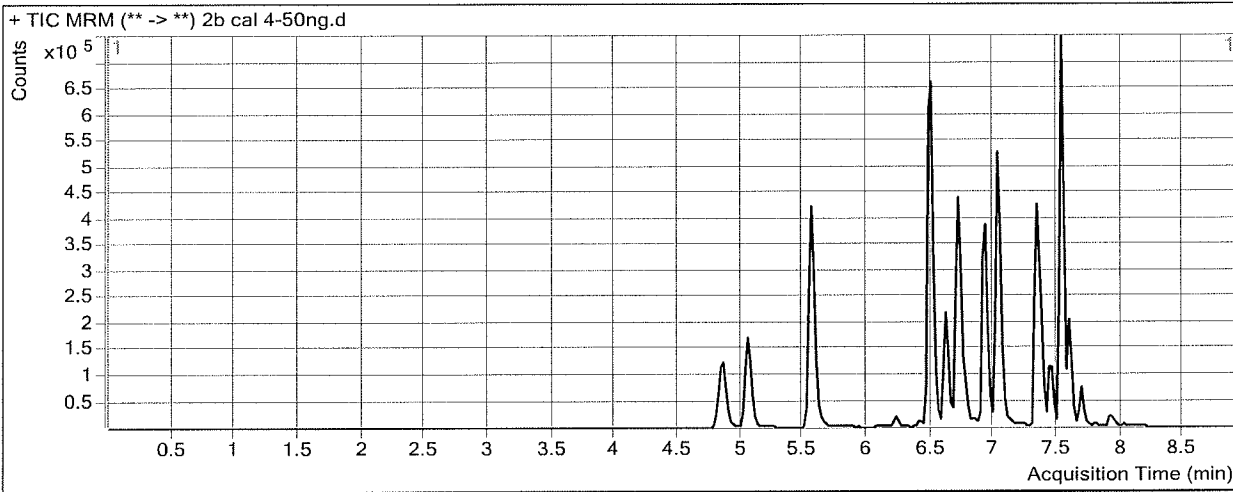
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 06:45 **Data File** 2b cal 4-50ng.d
Sample Type Calibration **Sample Name** 2b cal 4-50ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-D5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 264847 | 144913 | 1.8276 | 54.0333 |
| MDMA | MDMA-D6 | 5.048 | 359862 | 43658 | 8.2427 | 53.0370 |
| MDEA | MDEA-D6 | 5.548 | 407978 | 457340 | 0.8921 | 51.9003 |
| Primidone | Tapentadol-D3 | 6.215 | 39951 | 905915 | 0.0441 | 46.8476 |
| Tapentadol | Tapentadol-D3 | 6.503 | 974120 | 905915 | 1.0753 | 51.0863 |
| Meperidine | Meperidine-D4 | 6.616 | 248543 | 253477 | 0.9805 | 49.3442 |
| Normeperidine | Normeperidine-D4 | 6.724 | 273973 | 187276 | 1.4629 | 52.6829 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 286508 | 257154 | 1.1141 | 52.1209 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 482704 | 330499 | 1.4605 | 49.7134 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 571272 | 585539 | 0.9756 | 50.9631 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 429546 | 57630 | 7.4535 | 46.7105 |
| Zaleplon | Zaleplon-D4 | 7.341 | 257341 | 188624 | 1.3643 | 48.8490 |
| Phenytoin | Phenytoin-D10 | 7.389 | 14695 | 15705 | 0.9357 | 46.8649 |
| Verapamil | Trimipramine-D3 | 7.435 | 241889 | 156679 | 1.5439 | 51.0992 |
| Amoxapine | Clomipramine-D3 | 7.500 | 30842 | 51923 | 0.5940 | 52.9552 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 399283 | 89724 | 4.4501 | 53.9537 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 249790 | 57630 | 4.3344 | 43.2441 |
| Imipramine | Imipramine-D3 | 7.528 | 176050 | 97603 | 1.8037 | 51.6889 |
| Desipramine | Desipramine-D3 | 7.549 | 113478 | 113164 | 1.0028 | 49.8255 |
| Trimipramine | Trimipramine-D3 | 7.586 | 158833 | 156679 | 1.0137 | 47.3563 |
| Clomipramine | Clomipramine-D3 | 7.681 | 54581 | 51923 | 1.0512 | 59.7118 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.908 | 29434 | 21511 | 1.3684 | 44.1790 |

ISP FORENSICS - Cd'A Instrument # 62340

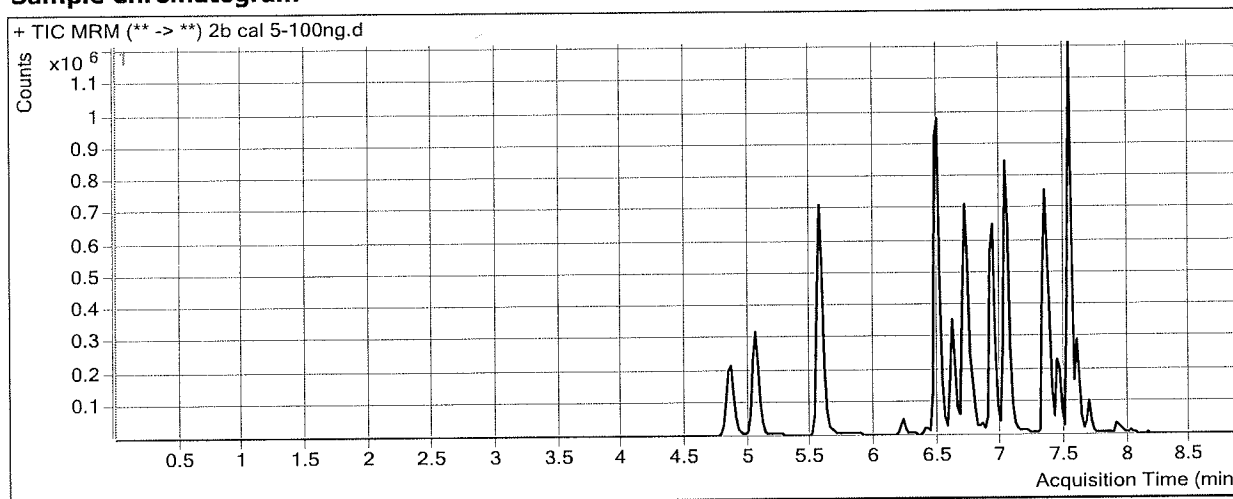
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 07:11 **Data File** 2b cal 5-100ng.d
Sample Type Calibration **Sample Name** 2b cal 5-100ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-E5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 511805 | 142651 | 3.5878 | 106.8065 |
| MDMA | MDMA-D6 | 5.048 | 730431 | 44196 | 16.5272 | 106.5269 |
| MDEA | MDEA-D6 | 5.548 | 792989 | 459047 | 1.7275 | 100.1499 |
| Primidone | Tapentadol-D3 | 6.215 | 82514 | 874749 | 0.0943 | 99.6821 |
| Tapentadol | Tapentadol-D3 | 6.503 | 1850377 | 874749 | 2.1153 | 100.4044 |
| Meperidine | Meperidine-D4 | 6.616 | 492119 | 248352 | 1.9815 | 99.1570 |
| Normeperidine | Normeperidine-D4 | 6.724 | 519913 | 175918 | 2.9554 | 106.9950 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 572083 | 257097 | 2.2252 | 104.3353 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 935386 | 319801 | 2.9249 | 99.5127 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 1115316 | 579040 | 1.9261 | 100.1484 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 880969 | 49195 | 17.9078 | 107.5059 |
| Zaleplon | Zaleplon-D4 | 7.341 | 496521 | 181196 | 2.7402 | 97.3659 |
| Phenytoin | Phenytoin-D10 | 7.389 | 32604 | 15921 | 2.0478 | 101.5689 |
| Verapamil | Trimipramine-D3 | 7.415 | 435378 | 130438 | 3.3378 | 111.5874 |
| Amoxapine | Clomipramine-D3 | 7.500 | 52600 | 41718 | 1.2608 | 112.5346 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 728610 | 89704 | 8.1223 | 98.4768 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 471687 | 49195 | 9.5882 | 93.4501 |
| Imipramine | Imipramine-D3 | 7.528 | 293607 | 85469 | 3.4352 | 97.4169 |
| Desipramine | Desipramine-D3 | 7.549 | 182994 | 93570 | 1.9557 | 96.8421 |
| Trimipramine | Trimipramine-D3 | 7.586 | 271784 | 130438 | 2.0836 | 96.6514 |
| Clomipramine | Clomipramine-D3 | 7.681 | 81372 | 41718 | 1.9505 | 111.6599 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 55674 | 18755 | 2.9685 | 93.1880 |

ISP FORENSICS - Cd'A Instrument # 62340

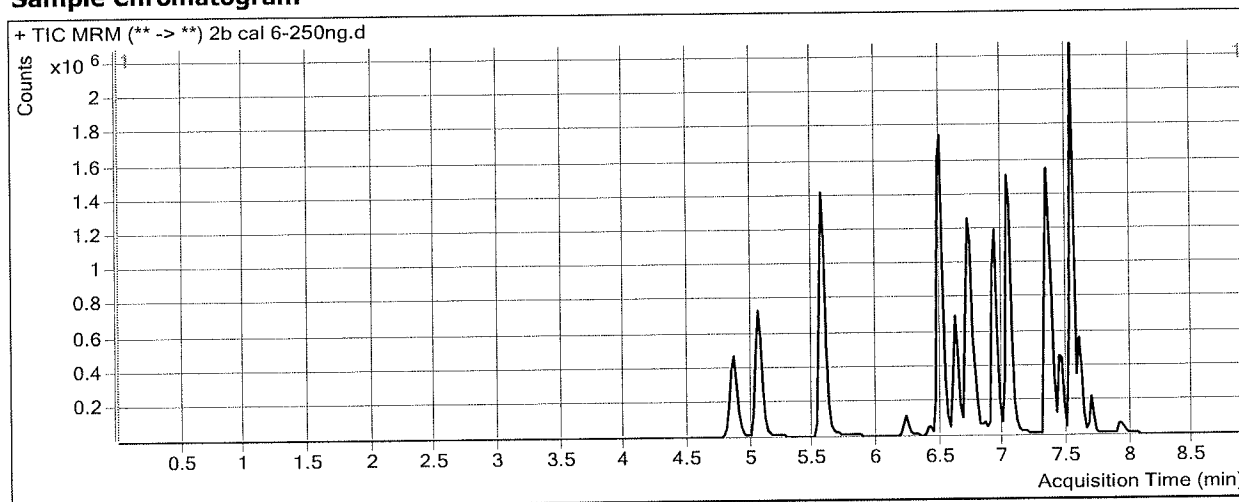
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 07:36 **Data File** 2b cal 6-250ng.d
Sample Type Calibration **Sample Name** 2b cal 6-250ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-F5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 1172172 | 142269 | 8.2391 | 246.2592 |
| MDMA | MDMA-D6 | 5.048 | 1671410 | 42985 | 38.8836 | 250.8759 |
| MDEA | MDEA-D6 | 5.548 | 1849030 | 440520 | 4.1974 | 242.8028 |
| Primidone | Tapentadol-D3 | 6.215 | 203671 | 764577 | 0.2664 | 280.6609 |
| Tapentadol | Tapentadol-D3 | 6.503 | 3966572 | 764577 | 5.1879 | 246.1064 |
| Meperidine | Meperidine-D4 | 6.616 | 1171188 | 236555 | 4.9510 | 246.9258 |
| Normeperidine | Normeperidine-D4 | 6.724 | 1038375 | 151637 | 6.8477 | 248.6367 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 1323306 | 245417 | 5.3921 | 253.1702 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 1945791 | 260596 | 7.4667 | 253.9667 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 2384975 | 519448 | 4.5914 | 238.0624 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 1926462 | 35427 | 54.3783 | 319.5947 |
| Zaleplon | Zaleplon-D4 | 7.341 | 1074299 | 153231 | 7.0110 | 247.9582 |
| Phenytoin | Phenytoin-D10 | 7.389 | 81463 | 16783 | 4.8539 | 239.5922 |
| Verapamil | Trimipramine-D3 | 7.415 | 961090 | 129513 | 7.4208 | 249.2556 |
| Amoxapine | Clomipramine-D3 | 7.500 | 117116 | 48467 | 2.4164 | 215.7770 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 1529850 | 75110 | 20.3681 | 246.9488 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 844158 | 35427 | 23.8281 | 229.5286 |
| Imipramine | Imipramine-D3 | 7.528 | 740344 | 82351 | 8.9901 | 253.1091 |
| Desipramine | Desipramine-D3 | 7.549 | 429692 | 85497 | 5.0258 | 248.3186 |
| Trimipramine | Trimipramine-D3 | 7.586 | 661178 | 129513 | 5.1051 | 235.8669 |
| Clomipramine | Clomipramine-D3 | 7.681 | 215679 | 48467 | 4.4500 | 256.0383 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.929 | 138111 | 18769 | 7.3584 | 227.6414 |

ISP FORENSICS - Cd'A Instrument # 62340

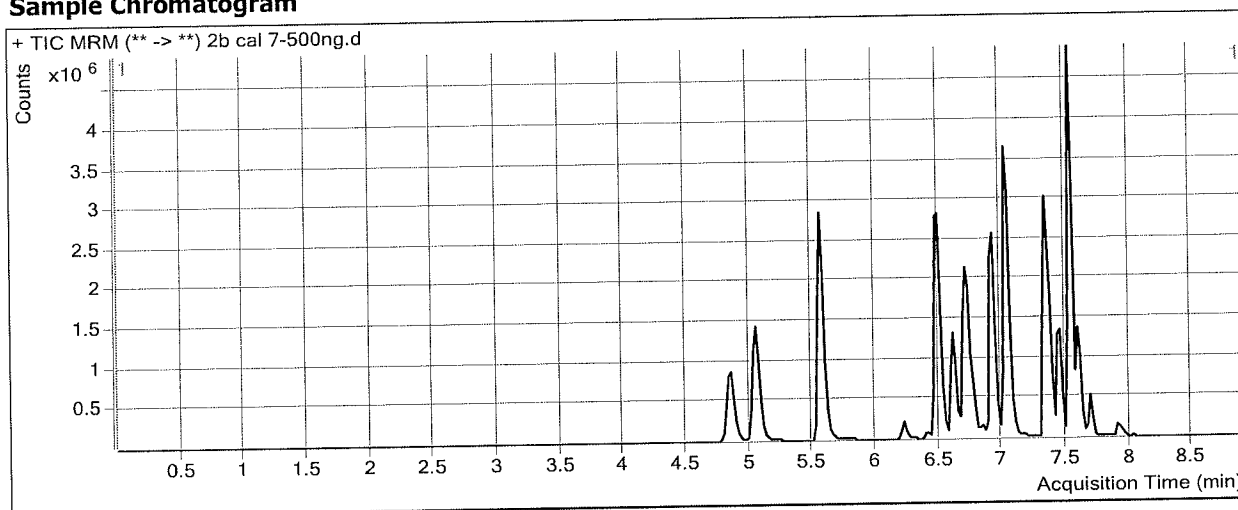
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 08:02 **Data File** 2b cal 7-500ng.d
Sample Type Calibration **Sample Name** 2b cal 7-500ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-G5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.850 | 2360239 | 137588 | 17.1544 | 513.5520 |
| MDMA | MDMA-D6 | 5.048 | 3419333 | 43668 | 78.3029 | 505.3941 |
| MDEA | MDEA-D6 | 5.548 | 3847155 | 440053 | 8.7425 | 505.3108 |
| Primidone | Tapentadol-D3 | 6.215 | 419469 | 704417 | 0.5955 | 626.8303 |
| Tapentadol | Tapentadol-D3 | 6.503 | 7420178 | 704417 | 10.5338 | 499.6053 |
| Meperidine | Meperidine-D4 | 6.616 | 2459353 | 243862 | 10.0850 | 502.4083 |
| Normeperidine | Normeperidine-D4 | 6.724 | 2069450 | 149338 | 13.8574 | 503.7200 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 2428604 | 230558 | 10.5336 | 494.8055 |
| Pentazocine | Pentazocine-13C3 | 6.919 | 4572736 | 306695 | 14.9097 | 507.0833 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 5946855 | 648543 | 9.1696 | 474.9660 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 4658225 | 21082 | 220.9623 | 1288.3409 |
| Zaleplon | Zaleplon-D4 | 7.341 | 1811222 | 130678 | 13.8602 | 489.4684 |
| Phenytoin | Phenytoin-D10 | 7.389 | 163754 | 15658 | 10.4582 | 515.2557 |
| Verapamil | Trimipramine-D3 | 7.435 | 2912104 | 169692 | 17.1611 | 577.6752 |
| Amoxapine | Clomipramine-D3 | 7.500 | 244242 | 61103 | 3.9972 | 357.0146 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 4197326 | 96338 | 43.5688 | 528.2427 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 1112217 | 21082 | 52.7579 | 505.9861 |
| Imipramine | Imipramine-D3 | 7.528 | 1774969 | 110541 | 16.0571 | 451.1834 |
| Desipramine | Desipramine-D3 | 7.549 | 990651 | 95481 | 10.3754 | 512.2634 |
| Trimipramine | Trimipramine-D3 | 7.586 | 1812650 | 169692 | 10.6820 | 492.8234 |
| Clomipramine | Clomipramine-D3 | 7.681 | 534980 | 61103 | 8.7554 | 504.7354 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 364212 | 23449 | 15.5322 | 477.9838 |

ISP FORENSICS - Cd'A Instrument # 62340

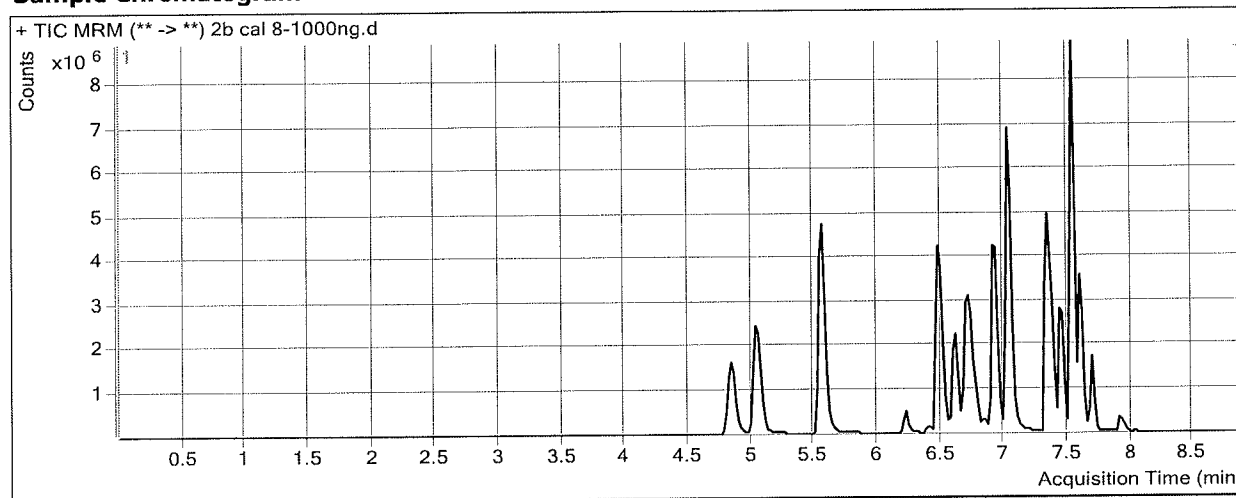
Multi-Drug Confirmatory Analysis Report

Batch Data Path D:\2017 Data\121917 mdq p1 and p2\QuantResults\121917 mdq p2b.batch.bin
Analysis Time 12/21/2017 12:27 PM **Analyst Name** ISP Tox
Report Time 12/21/2017 12:27 PM **Reporter Name** ISP Tox
Last Calib Update 12/21/2017 12:27 PM **Batch State** Processed

Analysis Info

Acq Time 2017-12-20 08:27 **Data File** 2b cal 8-1000ng.d
Sample Type Calibration **Sample Name** 2b cal 8-1000ng
Dilution 1 **Acq Method** MDQ Panel 2b 8-8-17.m
Position P1-H5 **Sample Info**
Inj Vol -1 **Comment** AM 28 Panel 2

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|-------------------------|-------|----------|-----------|------------|------------|
| MDA | MDA-D5 | 4.830 | 4453227 | 136132 | 32.7127 | 980.0102 |
| MDMA | MDMA-D6 | 5.028 | 6399516 | 41925 | 152.6436 | 985.3889 |
| MDEA | MDEA-D6 | 5.548 | 7302598 | 421522 | 17.3243 | 1000.9673 |
| Primidone | Tapentadol-D3 | 6.215 | 923910 | 561595 | 1.6452 | 1730.9461 |
| Tapentadol | Tapentadol-D3 | 6.483 | 11881677 | 561595 | 21.1570 | 1003.3561 |
| Meperidine | Meperidine-D4 | 6.616 | 4587528 | 227795 | 20.1388 | 1002.7103 |
| Normeperidine | Normeperidine-D4 | 6.703 | 3526626 | 129709 | 27.1888 | 988.8502 |
| 7-aminoflunitrazepam | 7-aminoflunitrazepam-D7 | 6.736 | 4190981 | 217091 | 19.3052 | 907.0452 |
| Pentazocine | Pentazocine-13C3 | 6.899 | 8195794 | 281524 | 29.1123 | 990.0713 |
| Phencyclidine | Phencyclidine-D5 | 7.028 | 12425558 | 620828 | 20.0145 | 1036.1468 |
| Flurazepam | Flunitrazepam-D7 | 7.335 | 9552032 | 12331 | 774.6629 | 4508.3097 |
| Zaleplon | Zaleplon-D4 | 7.341 | 2575559 | 89305 | 28.8402 | 1017.6818 |
| Phenytoin | Phenytoin-D10 | 7.389 | 364775 | 16382 | 22.2671 | 1096.1064 |
| Verapamil | Trimipramine-D3 | 7.415 | 6123957 | 226174 | 27.0763 | 911.9924 |
| Amoxapine | Clomipramine-D3 | 7.500 | 472149 | 109595 | 4.3081 | 384.7892 |
| Propoxyphene | Propoxyphene-D11 | 7.517 | 7944188 | 101972 | 77.9054 | 944.5523 |
| Flunitrazepam | Flunitrazepam-D7 | 7.520 | 1322956 | 12331 | 107.2908 | 1027.1107 |
| Imipramine | Imipramine-D3 | 7.528 | 4498434 | 111218 | 40.4469 | 1134.7842 |
| Desipramine | Desipramine-D3 | 7.549 | 2191086 | 108867 | 20.1262 | 993.3635 |
| Trimipramine | Trimipramine-D3 | 7.586 | 5041129 | 226174 | 22.2887 | 1027.6074 |
| Clomipramine | Clomipramine-D3 | 7.681 | 1842061 | 109595 | 16.8078 | 969.8707 |
| Norpropoxyphene | Norpropoxyphene-D5 | 7.909 | 808653 | 23462 | 34.4667 | 1057.9036 |