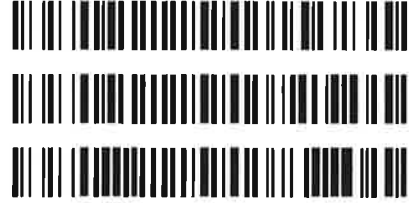


TS

6/18/2018

Worklist: 2495

| <u>LAB CASE</u> | <u>ITEM</u> | <u>TASK ID</u> | <u>DESCRIPTION</u> |
|-----------------|-------------|----------------|---|
| M2018-2479 | 1 | 118877 | AM 28 Blood Multi-Drug Quant Panel 2 by LC-QC |
| M2018-2617 | 2 | 118895 | AM 28 Blood Multi-Drug Quant Panel 2 by LC-QC |
| P2018-1522 | 1 | 118896 | AM 28 Blood Multi-Drug Quant Panel 2 by LC-QC |



AM# 28: Multi-Drug Quantitation in Blood by LC-MS/MS

TS

Extraction Date: 06/18/2018
Plate lot#: 0530563

Analyst: Tamara Salazar
Plate Expiration: 05/21/2019

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

Mobile phase B: 0.01% Formic Acid in MeOH
Ethyl Acetate 20% Methanol in Water

Blank Blood Lot: 361331-1

Column: Agilent 120 EC-C18 (2.1x 100-4um)

LCMS-QQQ ID: 59740

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: 3** in wells of analytical (standards) plate.
- 3. Place on shaking incubator at ambient temp., 900rpm for 15 minutes. *Shaker ID: 067105*
- 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: 067104
- 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: 067103
- 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: *C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1and P2 TS*
Batch Name: *p2a MDQ 61818 final*
p2b MDQ 61818 final
- 2. Make necessary changes to integration limits
- 3. Integration linear and R² 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% (tramadol RT +/-2%), S/N for primary transition >10 and secondary transitions >5.
- 5. Did all QCs pass for each analyte? Y / N _____ Add Control data to QC tracking spreadsheet.
- 6. Central File Packet to include: LIMS Worklist, Method Checklist, Calibration and Control Reports.

AM# 28: Multi-Drug Quantitation in Blood by LC-MS/MS Cont.

5

Extraction Date: 06/18/2018

Plate lot#: 0530563

Analyst: Tamara Salazar

Plate Expiration: 05/21/2019

COMMENTS:

P2a evaluated for: Pseudoephedrine

Zopiclone

Amitriptyline

Nortriptyline

Midazolam

Alpha-hydroxymidazolam

Chlordiazepoxide

P1b evaluated for: Flunitrazepam

Curves Limited: Nortriptyline 5-500



Idaho State Police Forensic Services

TS

AM #28 Blood Multi-Drug Confirmatory Analysis by LCMS-QQQ

Analyst: Tamara Salazar
Extraction Date: 06/18/18
Worklist Number: 2495

| Reagent | Lot Number | Expiration Date | Date in Service | Date Out of Service | Initials |
|--|--|-----------------|--|---------------------|----------|
| ToxBox Benzo/Opiate Quant Plate | 0530563 | 05/21/19 | 06/05/18 | | |
| Negative Blood | 361331-1 | | 05/25/18 | | |
| Methanol External Control Solution | 022018 | 02/20/19 | 02/20/18 | | |
| Blood External Control Solution | WS061218 | 06/12/19 | 06/12/18 | | |
| 0.5M Ammonium Hydroxide | 032918 | 09/29/18 | 03/29/18 | | |
| Formic Acid (LCMS Grade) | 095180B | | 11/15/17 | | |
| Water (LCMS Grade) | 177528 | | 03/05/18 | | |
| Methanol (LCMS Grade) | 177145 | | 04/11/18 | | |
| Ammonium Formate (LCMS Grade) | A0287238 | | 01/30/17 | | |
| Ammonium Hydroxide (ACS grade or better) | 053200 | | 08/03/07 | | |
| Ethyl Acetate (ACS grade or better) | 100362 | | 06/26/17 | | |
| 20% LCMS Methanol in LCMS Water | 052918 06/18/18 ↗ | Make Fresh | 052918 06/18/18 ↗ | | |
| 5mM Ammonium Formate + 0.01% Formic Acid in Water (Mobile Phase A) | 051018 | | 05/10/18 | | |
| 0.01% Formic Acid in Methanol (Mobile Phase B) | 061818 | | 06/18/18 | | |
| Needle Rinse--75% LCMS MeOH in LCMS Water | 061818 | | 06/18/18 | | |

Methanol External Control Solution (Lot: 022018)
100 ul each 1 mg/mL stock solution in 9600 ul MeOH

| Component | Source | Source Lot Number | Expiration Date |
|------------------|----------------|-------------------|-----------------|
| Methanol (LCMS) | Fisher | 172516 | |
| Zopiclone | Cerilliant | FE08291601 | 12/31/2021 |
| Chlordiazepoxide | Cerilliant | FE07241502 | 08/31/2020 |
| Flunitrazepam | Cerilliant | FE08051602 | 08/31/2021 |
| Prepared: | 02/20/18 | | |
| Prepared By: | Tamara Salazar | | |
| Expires: | 02/20/19 | | |

TS

AM #28: Blood Multi-Drug Confirmatory Analysis by LCMS-QQQ

Blood External Control Solution (Lot: WS061218)

50 ul of methanol external control solution was added to 9950 ul of blood.

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|------------------------------------|----------------|--------------------------|
| Negative Blood | Hemostat | 361331-1 |
| Methanol External Control Solution | | 121217 |
| Prepared: | 06/12/18 | |
| Prepared by: | Tamara Salazar | |
| Expires: | 06/12/19 | |

0.5M Ammonium Hydroxide (Lot: 032918)

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|--|----------------|--------------------------|
| Ammonium Hydroxide (ACS grade or better) | Fisher | 053200 |
| DI Water | - | - |
| Prepared: | 03/29/18 | |
| Prepared By: | Tamara Salazar | |
| Expires: | 09/29/18 | |

20% LCMS Methanol in LCMS Water (Lot: Prepared Fresh)

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|--------------------|---------------|--------------------------|
| MeOH (LCMS Grade) | Fisher | 177145 |
| Water (LCMS Grade) | Fisher | 177528 |

5mM Ammonium Formate + 0.01% Formic Acid in Water (Lot: 051018)

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|------------------|---------------|--------------------------|
| Ammonium Formate | Arcos Organic | A0287238 |
| LCMS Water | Fisher | 177528 |
| Formic Acid | Fisher | 095180B |
| Prepared: | 05/10/18 | |
| Prepared By: | Sarah Pickle | |

0.01% Formic Acid in LCMS Methanol (Mobile Phase B) (Lot: 031918)

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|-------------------|----------------|--------------------------|
| Formic Acid | Fisher | 095180B |
| MeOH (LCMS Grade) | Fisher | 172517 |
| Prepared: | 06/18/18 | |
| Prepared By: | Tamara Salazar | |

Needle Rinse (75% LCMS MeOH in LCMS Water) (Lot: 052918)

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|--------------------|----------------|--------------------------|
| MeOH (LCMS Grade) | Fisher | 177145 |
| Water (LCMS Grade) | Fisher | 177528 |
| Prepared: | 06/18/18 | |
| Prepared By: | Tamara Salazar | |

TS

ISP FORENSICS - Pocatello Instrument # 59740

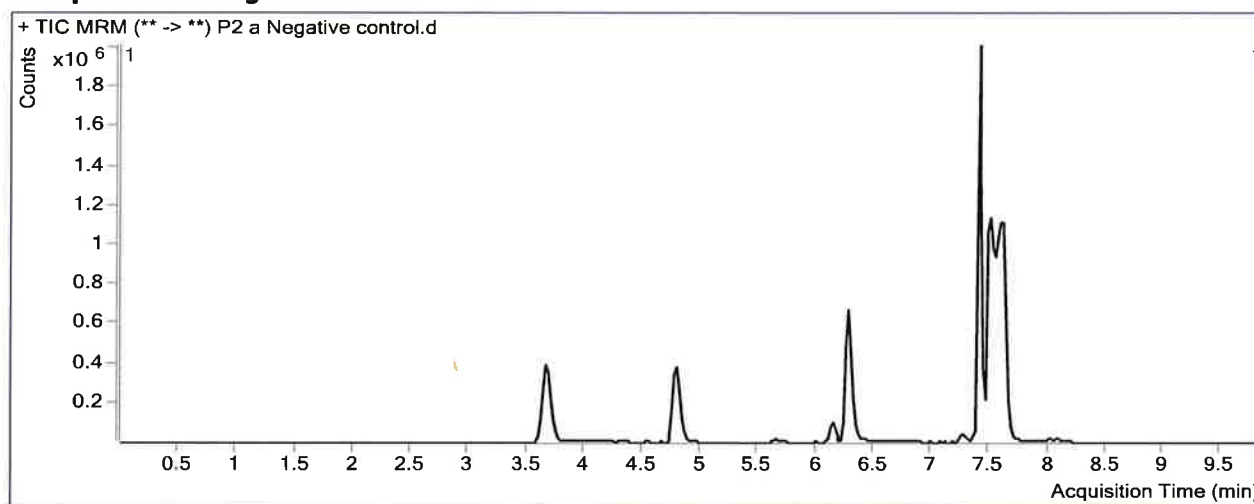
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 19:38 **Data File** P2 a Negative control.d
Sample Type Sample **Sample Name** P2 a Negative control.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 |
|-----------------|--------------------|-------|----------|-----------|------------|---------------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 4.161 | 13317 | 1948119 | 0.0068 | 2.0879 <i>LS TD</i> |

TS

ISP FORENSICS - Pocatello Instrument # 59740

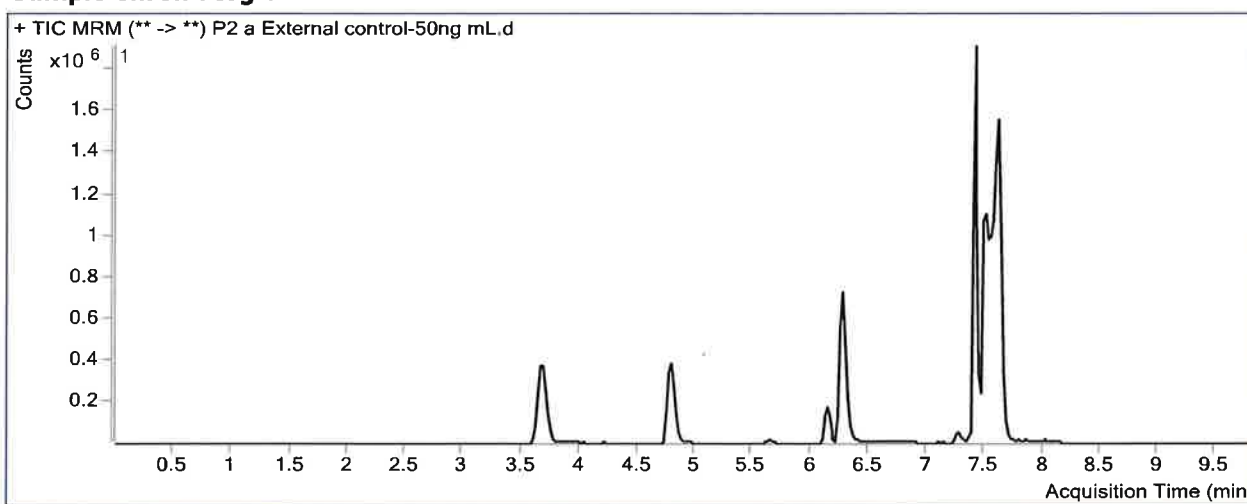
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 20:03 **Data File** P2 a External control-50ng mL.d
Sample Type Sample **Sample Name** P2 a External control-50ng mL.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.m
Position P1-F6 **Sample Info**
Inj Vol -1 **Comment** AM 28; WS061218 50 ng flunitrazepam, zopiclone, chlorodiazepoxide

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|-------------------------|---------------------|-------|----------|-----------|------------|----------------|
| <u>Zopiclone</u> | Zopiclone-D4 | 6.131 | 291411 | 243229 | 1.1981 | <u>39.9816</u> |
| <u>Chlordiazepoxide</u> | Chlordiazepoxide-D5 | 7.605 | 606155 | 2422794 | 0.2502 | <u>43.5703</u> |

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ISP FORENSICS - Pocatello Instrument # 59740

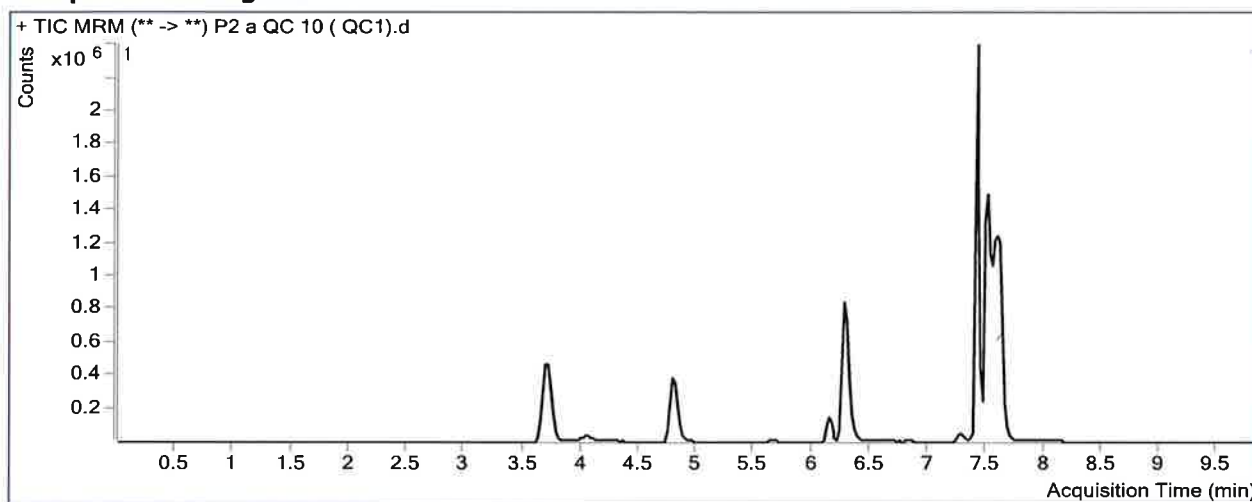
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 20:54 **Data File** P2 a QC 10 (QC1).d
Sample Type QC **Sample Name** P2 a QC 10 (QC1).d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.716 | 428838 | 1985573 | 0.2160 | 10.4420 |
| Zopiclone | Zopiclone-D4 | 6.151 | 100509 | 408780 | 0.2459 | 10.2461 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 17780 | 87121 | 0.2041 | 9.4148 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 12242 | 77052 | 0.1589 | 9.8091 |
| Midazolam | Midazolam-D4 | 7.543 | 87807 | 1960602 | 0.0448 | 9.6566 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 228595 | 1882820 | 0.1214 | 10.6827 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 131744 | 2431883 | 0.0542 | 10.9163 |

15

ISP FORENSICS - Pocatello Instrument # 59740

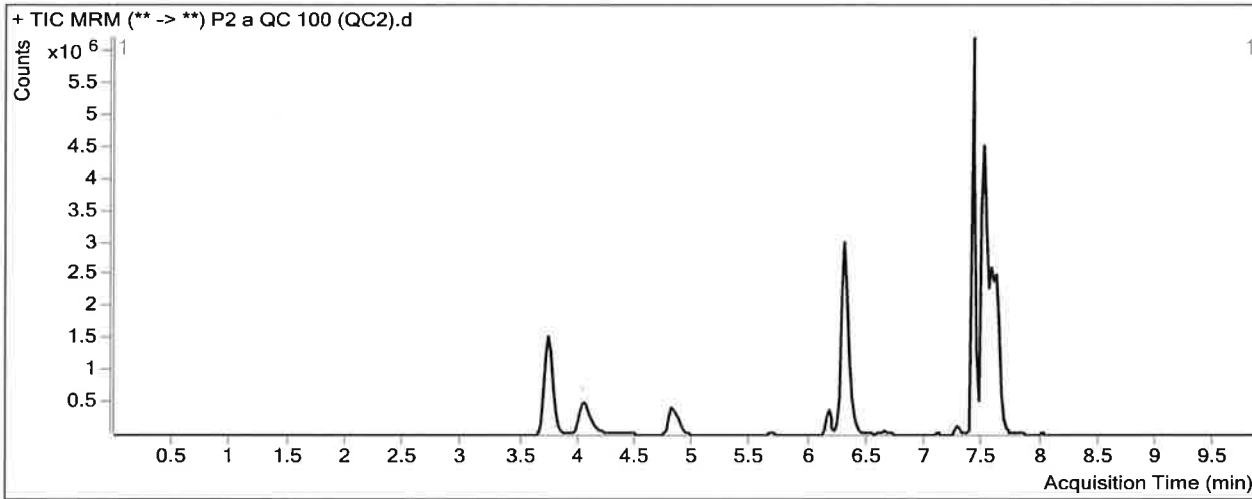
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 22:11 **Data File** P2 a QC 100 (QC2).d
Sample Type QC **Sample Name** P2 a QC 100 (QC2).d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 4876923 | 2023181 | 2.4105 | 98.1034 |
| Zopiclone | Zopiclone-D4 | 6.151 | 804268 | 273215 | 2.9437 | 94.4934 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 140002 | 64823 | 2.1598 | 89.5431 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 103690 | 51134 | 2.0278 | 110.7697 |
| Midazolam | Midazolam-D4 | 7.543 | 1062325 | 1873837 | 0.5669 | 99.6804 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.582 | 2255010 | 1625534 | 1.3872 | 111.3540 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 1219019 | 2093667 | 0.5822 | 98.8868 |

15

ISP FORENSICS - Pocatello Instrument # 59740

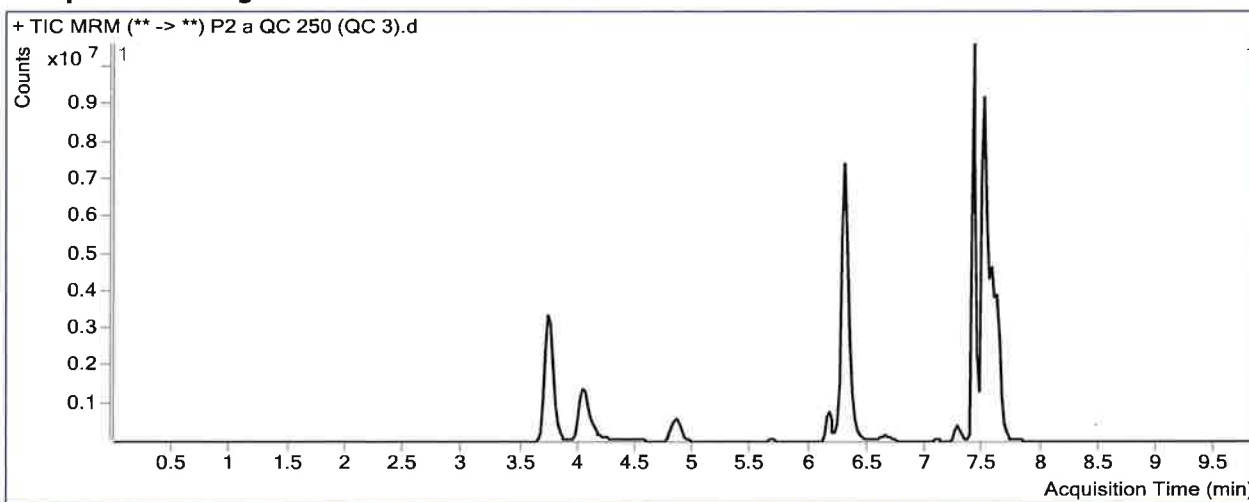
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:40 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 23:27 **Data File** P2 a QC 250 (QC 3).d
Sample Type QC **Sample Name** P2 a QC 250 (QC 3).d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 13265199 | 2081658 | 6.3724 | 256.3619 |
| Zopiclone | Zopiclone-D4 | 6.151 | 1978014 | 250192 | 7.9060 | 249.4540 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 681743 | 110283 | 6.1817 | 254.3314 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 503210 | 98623 | 5.1023 | 276.8600 |
| Midazolam | Midazolam-D4 | 7.543 | 2670597 | 1898433 | 1.4067 | 244.4753 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 4665848 | 1370409 | 3.4047 | 271.8028 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 2393961 | 1639573 | 1.4601 | 245.1312 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

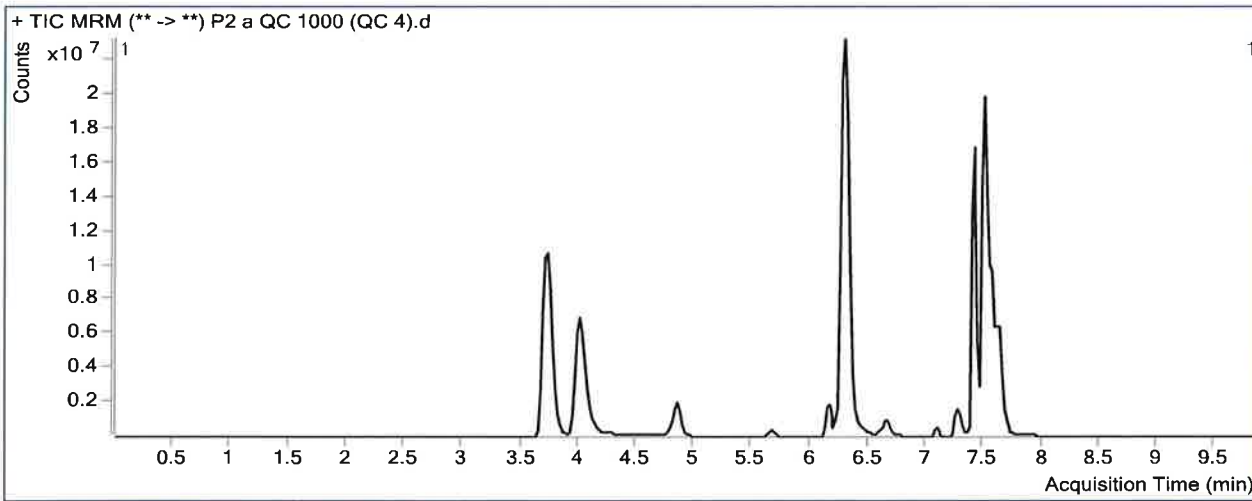
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:40 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-19 00:43 **Data File** P2 a QC 1000 (QC 4).d
Sample Type QC **Sample Name** P2 a QC 1000 (QC 4).d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 52369283 | 2015474 | 25.9836 | 1039.7331 |
| Zopiclone | Zopiclone-D4 | 6.151 | 5000080 | 151121 | 33.0865 | 1035.7844 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 1808060 | 79214 | 22.8250 | 936.2400 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 986552 | 44942 | 21.9515 | 1187.0690 |
| Midazolam | Midazolam-D4 | 7.543 | 8318913 | 1390246 | 5.9838 | 1033.6167 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.582 | 9285664 | 651552 | 14.2516 | 1134.4535 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.625 | 5465333 | 862424 | 6.3372 | 1057.5995 |

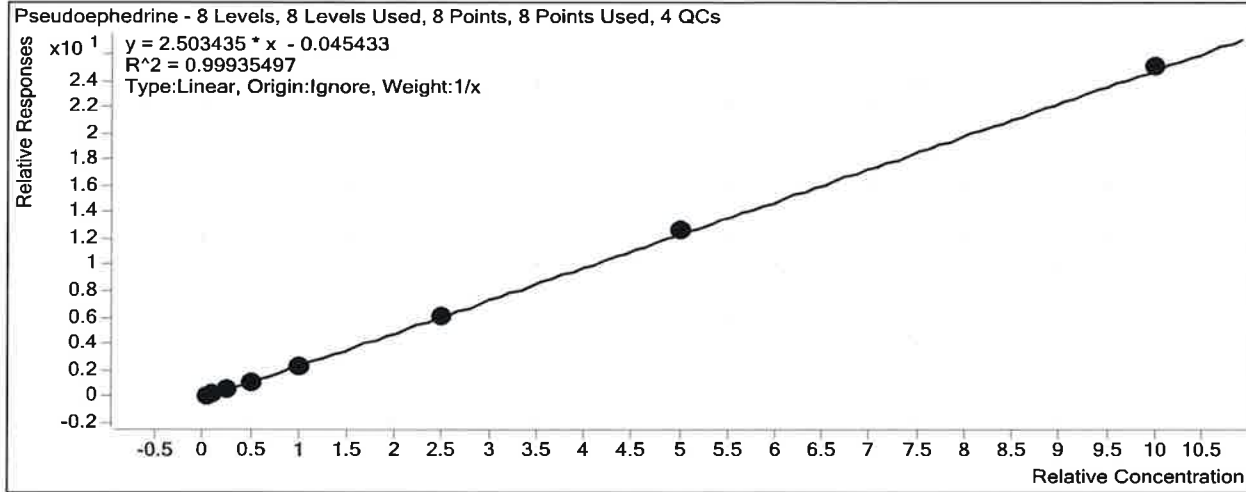
Outside curve range. TS

ISP Forensics Calibration Curve Report

TS

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

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



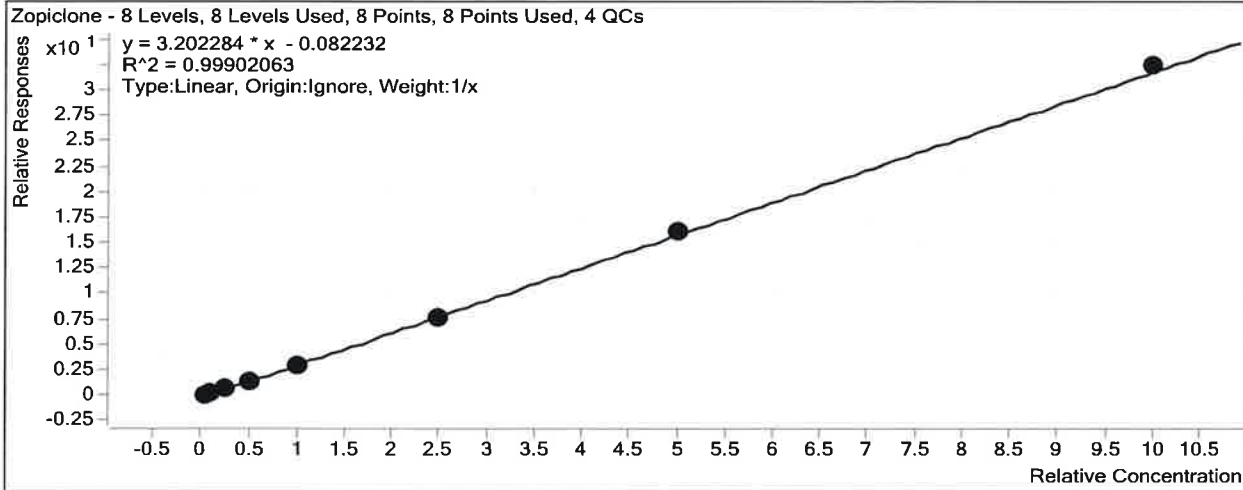
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 5.8 | 116.3 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.9 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 10.4 | 104.4 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 24.1 | 96.4 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 47.4 | 94.8 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 93.3 | 93.3 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 98.1 | 98.1 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 244.3 | 97.7 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 256.4 | 102.5 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 511.0 | 102.2 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1004.3 | 100.4 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1039.7 | 104.0 |

ISP Forensics Calibration Curve Report

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Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

Target Compound Zopiclone
Internal Standard Zopiclone-D4



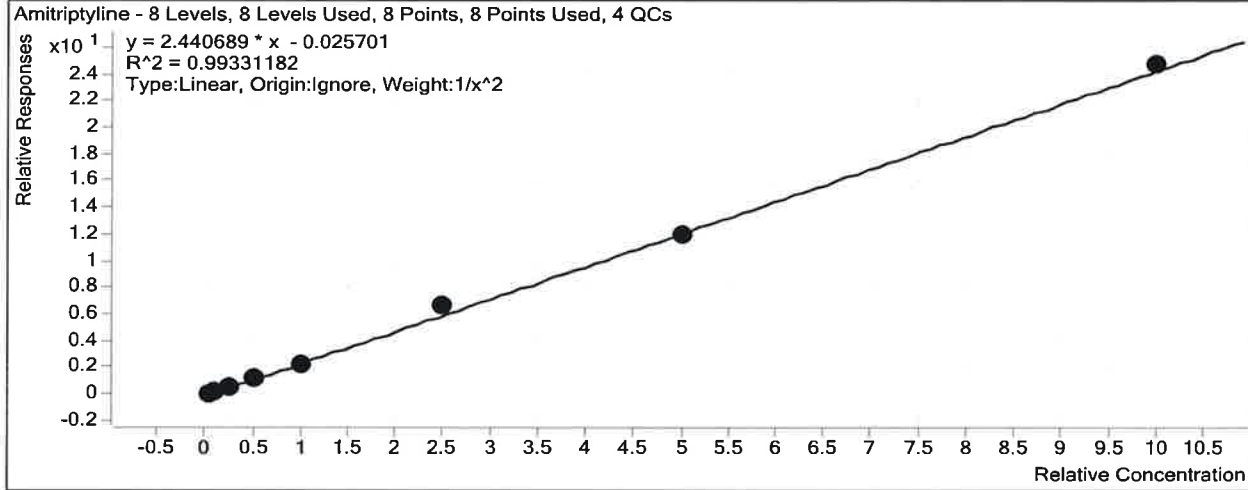
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 6.0 | 119.6 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 10.1 | 101.4 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.5 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 23.5 | 94.1 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 46.9 | 93.8 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 91.9 | 91.9 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 94.5 | 94.5 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 240.9 | 96.4 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 249.5 | 99.8 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 506.9 | 101.4 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1013.7 | 101.4 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1035.8 | 103.6 |

ISP Forensics Calibration Curve Report

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Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

Target Compound Amitriptyline
Internal Standard Amitriptyline-D3



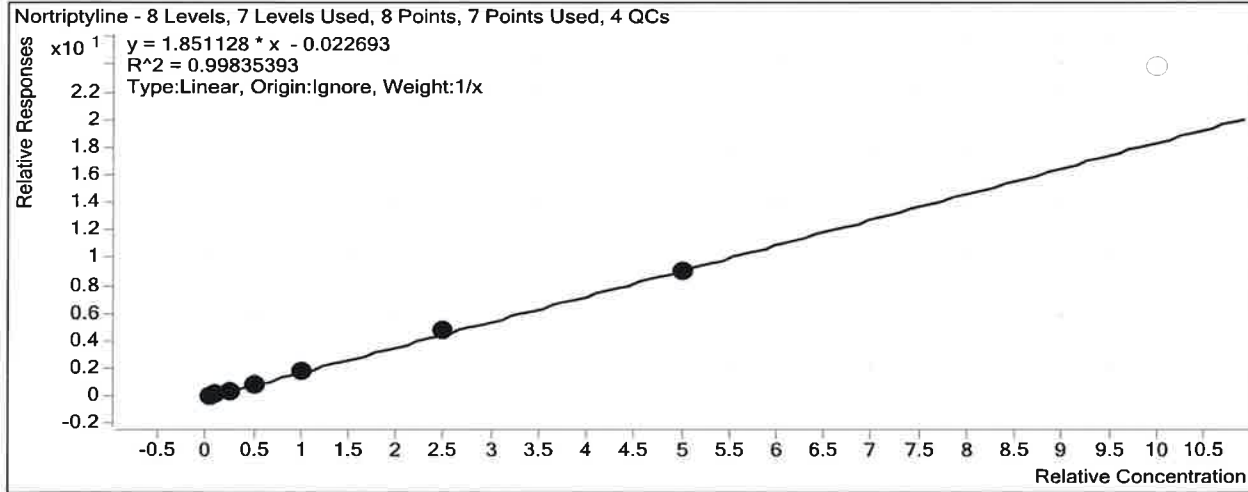
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 5.3 | 106.5 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 8.6 | 85.9 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 9.4 | 94.1 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 25.9 | 103.6 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 49.3 | 98.5 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 96.0 | 96.0 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 89.5 | 89.5 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 272.4 | 109.0 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 254.3 | 101.7 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 495.1 | 99.0 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1014.0 | 101.4 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 936.2 | 93.6 |

ISP Forensics Calibration Curve Report

TS

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

Target Compound Nortriptyline
Internal Standard Nortriptyline-D3



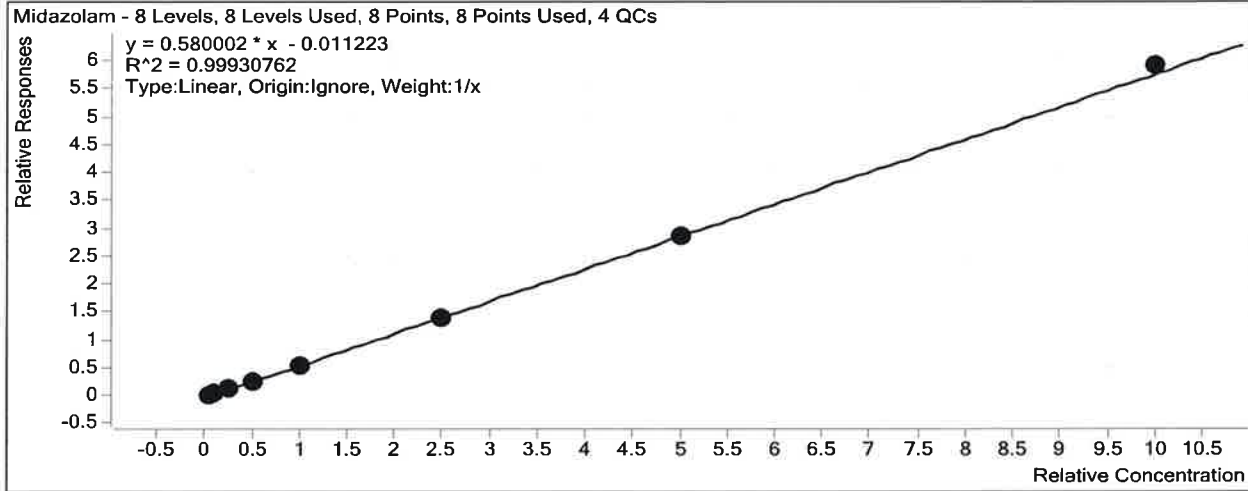
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 5.5 | 109.3 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 9.7 | 96.6 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 9.8 | 98.1 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 23.9 | 95.8 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 46.3 | 92.6 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 102.8 | 102.8 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 110.8 | 110.8 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 262.7 | 105.1 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 276.9 | 110.7 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 489.1 | 97.8 |
| P2 a cal 8-1000ng.d | 8 | <input type="checkbox"/> | 1000 | 1288.8 | 128.9 |
| P2 a QC 1000 (QC 4).d | 8 | <input type="checkbox"/> | 1000 | 1187.1 | 118.7 |

ISP Forensics Calibration Curve Report

TS

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

Target Compound Midazolam
Internal Standard Midazolam-D4



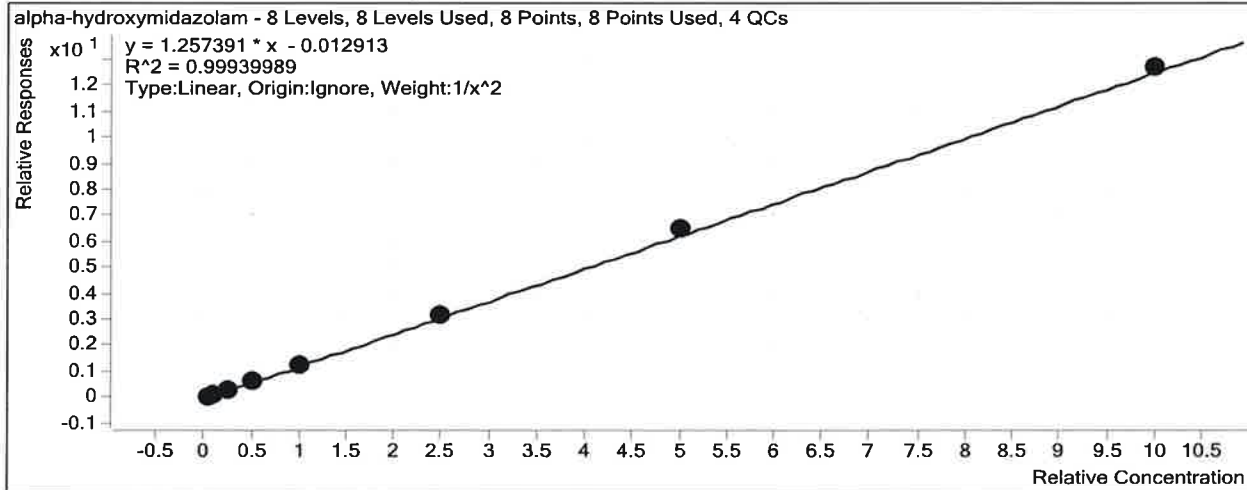
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 5.9 | 117.5 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.7 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 9.7 | 96.6 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 23.8 | 95.0 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 48.0 | 96.0 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 94.5 | 94.5 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 99.7 | 99.7 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 242.9 | 97.1 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 244.5 | 97.8 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 495.4 | 99.1 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1019.7 | 102.0 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1033.6 | 103.4 |

ISP Forensics Calibration Curve Report

B

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

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



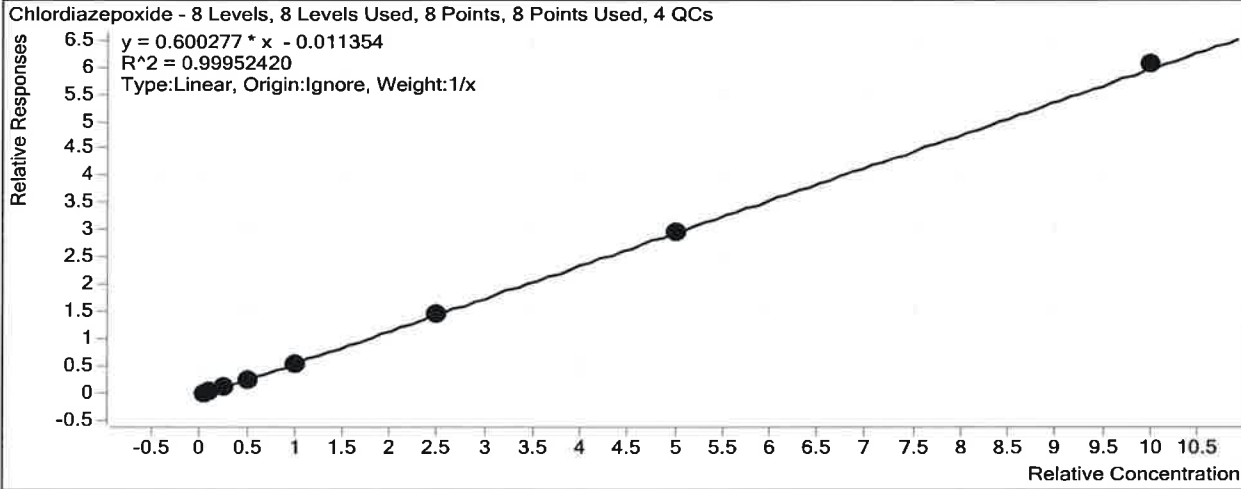
| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 5.1 | 101.0 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 9.9 | 98.7 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 10.7 | 106.8 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 24.9 | 99.4 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 49.2 | 98.4 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 97.0 | 97.0 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 111.4 | 111.4 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 251.4 | 100.5 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 271.8 | 108.7 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 519.5 | 103.9 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1010.8 | 101.1 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1134.5 | 113.4 |

ISP Forensics Calibration Curve Report

B

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:37 PM **Analyst Name** ISP TOX

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



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|-----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 a cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 6.0 | 119.2 |
| P2 a cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 9.0 | 90.3 |
| P2 a QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 10.9 | 109.2 |
| P2 a cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 25.2 | 100.9 |
| P2 a cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 47.5 | 95.0 |
| P2 a cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 94.2 | 94.2 |
| P2 a QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 98.9 | 98.9 |
| P2 a cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 248.9 | 99.6 |
| P2 a QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 245.1 | 98.1 |
| P2 a cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 499.5 | 99.9 |
| P2 a cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1009.7 | 101.0 |
| P2 a QC 1000 (QC 4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1057.6 | 105.8 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

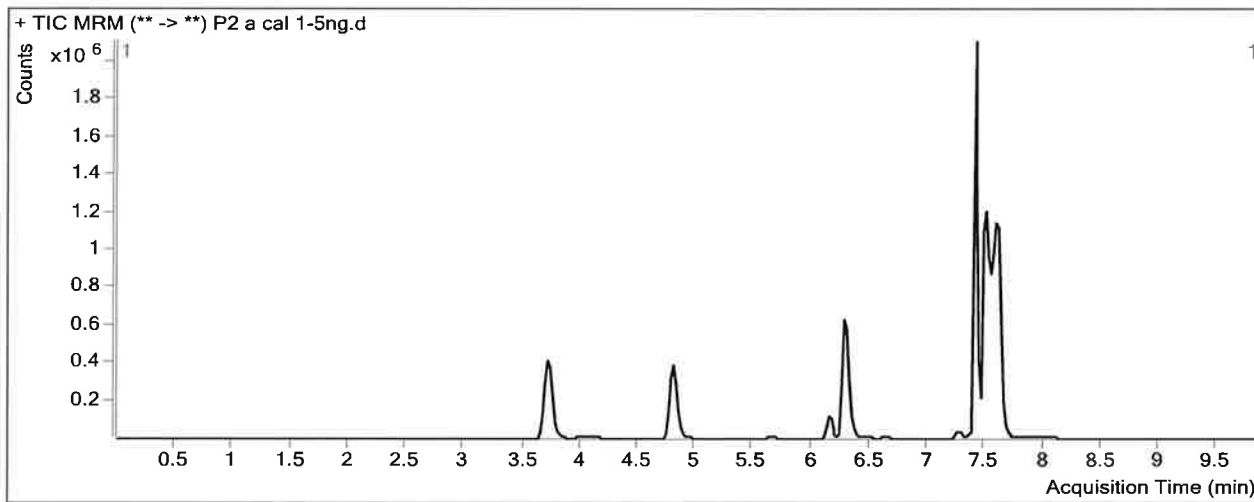
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:38 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 16:01 **Data File** P2 a cal 1-5ng.d
Sample Type Calibration **Sample Name** P2 a cal 1-5ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 182934 | 1827722 | 0.1001 | 5.8129 |
| Zopiclone | Zopiclone-D4 | 6.151 | 42394 | 388113 | 0.1092 | 5.9790 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 9191 | 88172 | 0.1042 | 5.3240 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 6150 | 78342 | 0.0785 | 5.4666 |
| Midazolam | Midazolam-D4 | 7.543 | 35829 | 1568459 | 0.0228 | 5.8735 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.582 | 95612 | 1889197 | 0.0506 | 5.0520 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 59765 | 2447284 | 0.0244 | 5.9597 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

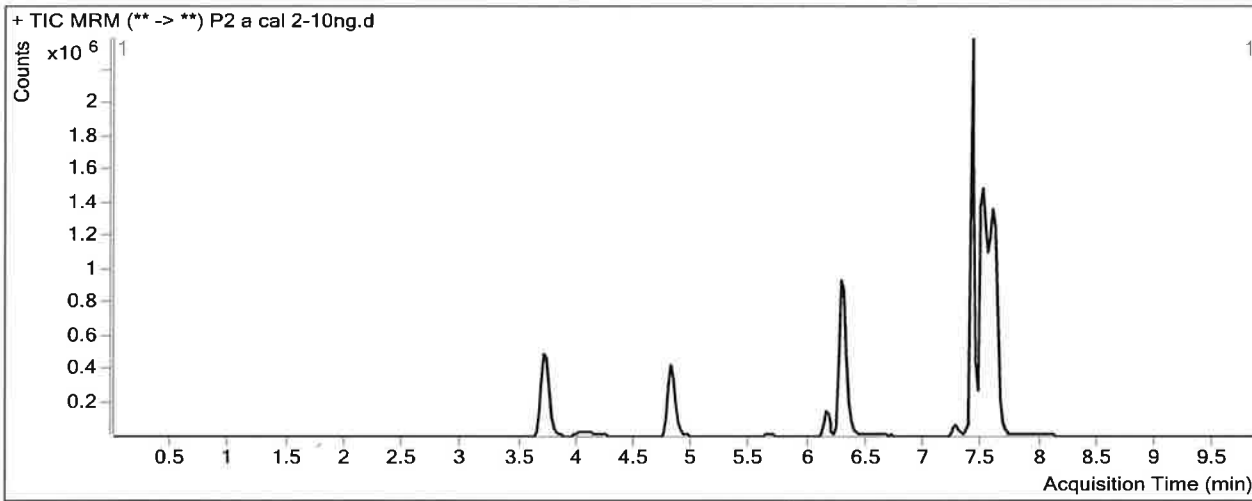
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:38 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 16:26 **Data File** P2 a cal 2-10ng.d
Sample Type Calibration **Sample Name** P2 a cal 2-10ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 412354 | 2039722 | 0.2022 | 9.8902 |
| Zopiclone | Zopiclone-D4 | 6.151 | 103767 | 427745 | 0.2426 | 10.1435 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 23615 | 128349 | 0.1840 | 8.5915 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 19948 | 127804 | 0.1561 | 9.6576 |
| Midazolam | Midazolam-D4 | 7.543 | 89967 | 1953698 | 0.0460 | 9.8745 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.582 | 227222 | 2044022 | 0.1112 | 9.8678 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 114562 | 2674451 | 0.0428 | 9.0275 |

B

ISP FORENSICS - Pocatello Instrument # 59740

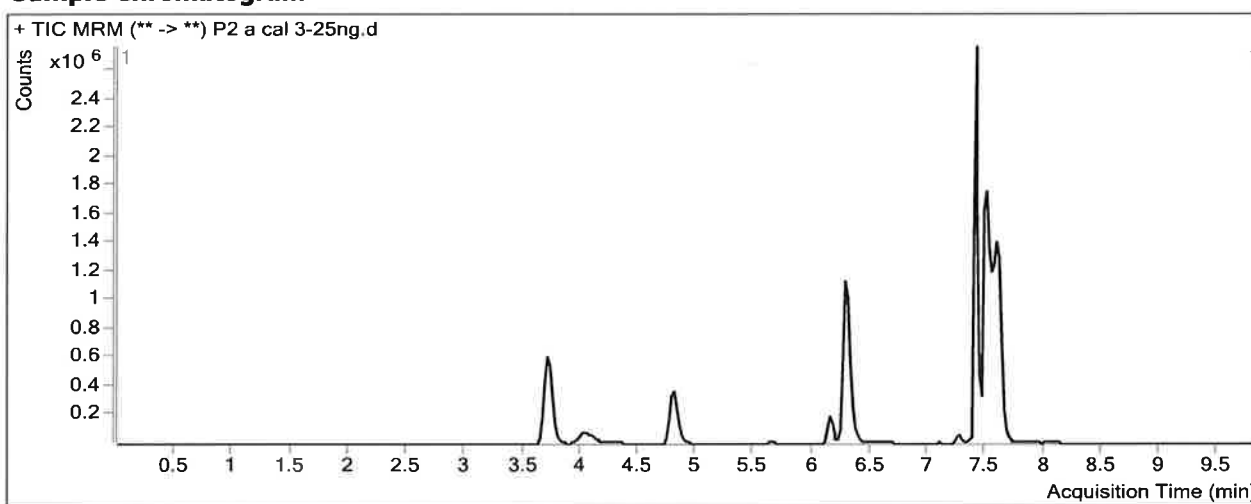
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:38 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 16:52 **Data File** P2 a cal 3-25ng.d
Sample Type Calibration **Sample Name** P2 a cal 3-25ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 1021425 | 1830505 | 0.5580 | 24.1043 |
| Zopiclone | Zopiclone-D4 | 6.131 | 233899 | 348419 | 0.6713 | 23.5316 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 33257 | 54813 | 0.6067 | 25.9124 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 20512 | 48772 | 0.4206 | 23.9462 |
| Midazolam | Midazolam-D4 | 7.543 | 219434 | 1733996 | 0.1265 | 23.7536 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 522161 | 1743103 | 0.2996 | 24.8508 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 311726 | 2224768 | 0.1401 | 25.2334 |

15

ISP FORENSICS - Pocatello Instrument # 59740

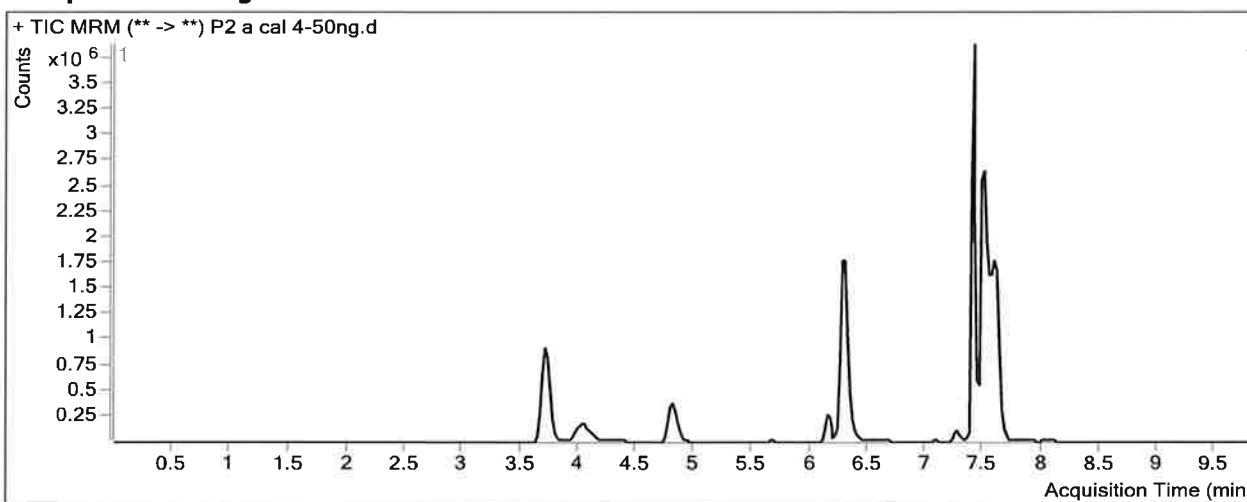
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:38 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 17:17 **Data File** P2 a cal 4-50ng.d
Sample Type Calibration **Sample Name** P2 a cal 4-50ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.736 | 2216194 | 1941295 | 1.1416 | 47.4164 |
| Zopiclone | Zopiclone-D4 | 6.131 | 473711 | 333652 | 1.4198 | 46.9042 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 121771 | 103481 | 1.1768 | 49.2669 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 74429 | 89198 | 0.8344 | 46.3025 |
| Midazolam | Midazolam-D4 | 7.543 | 487473 | 1823761 | 0.2673 | 48.0193 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 1059744 | 1750146 | 0.6055 | 49.1836 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 614524 | 2245617 | 0.2737 | 47.4796 |

B

ISP FORENSICS - Pocatello Instrument # 59740

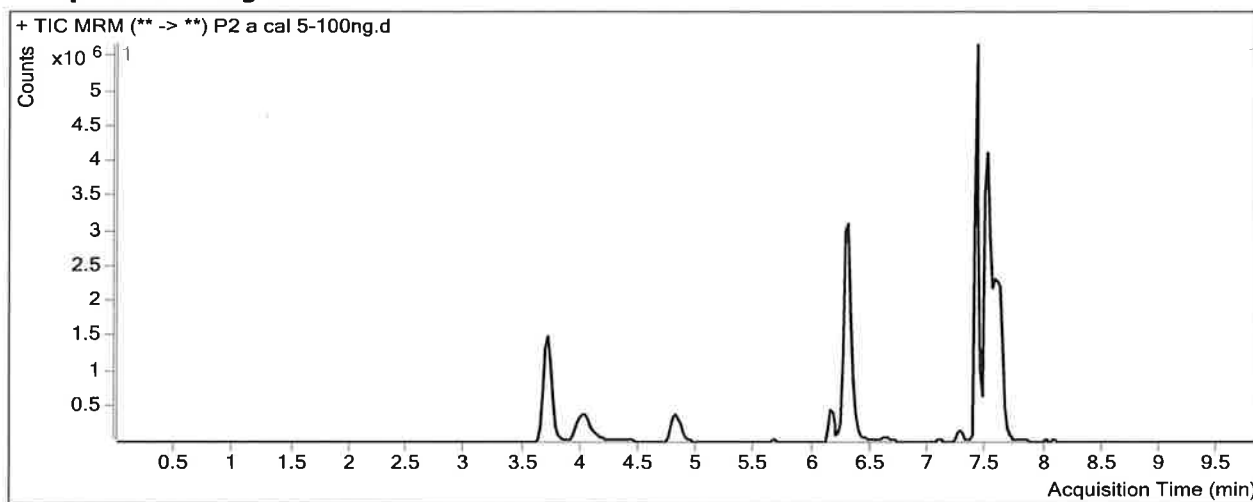
Multi-Drug Confirmatory Analysis Report

| | | | |
|--------------------------|---|----------------------|-----------|
| Batch Data Path | C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir | | |
| Analysis Time | 6/22/2018 1:37 PM | Analyst Name | ISPUser |
| Report Time | 6/22/2018 1:38 PM | Reporter Name | ISPUser |
| Last Calib Update | 6/22/2018 1:37 PM | Batch State | Processed |

Analysis Info

| | | | |
|--------------------|------------------|--------------------|------------------------|
| Acq Time | 2018-06-18 17:43 | Data File | P2 a cal 5-100ng.d |
| Sample Type | Calibration | Sample Name | P2 a cal 5-100ng.d |
| Dilution | 1 | Acq Method | MDQ Panel 2a 6-18-18.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 |
|------------------------|---------------------------|-------|----------|-----------|------------|------------|
| Pseudoephedrine | Pseudoephedrine-D3 | 3.716 | 4538418 | 1982462 | 2.2893 | 93.2605 |
| Zopiclone | Zopiclone-D4 | 6.131 | 994566 | 347515 | 2.8619 | 91.9396 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 189857 | 81898 | 2.3182 | 96.0353 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 127428 | 67777 | 1.8801 | 102.7912 |
| Midazolam | Midazolam-D4 | 7.543 | 978939 | 1822642 | 0.5371 | 94.5380 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 1967449 | 1630646 | 1.2065 | 96.9832 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 1123786 | 2027819 | 0.5542 | 94.2129 |

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ISP FORENSICS - Pocatello Instrument # 59740

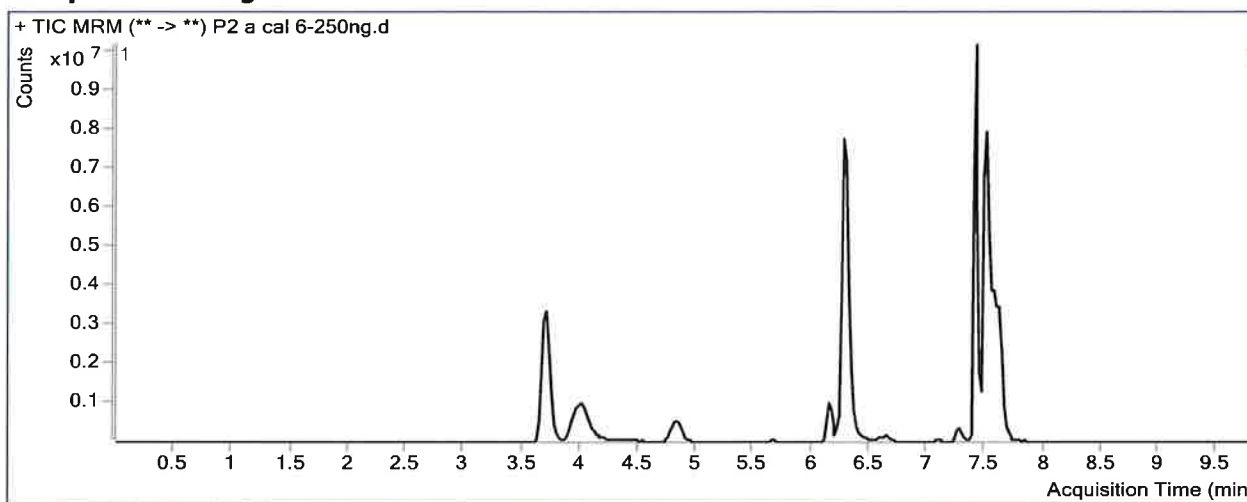
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:38 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 18:09 **Data File** P2 a cal 6-250ng.d
Sample Type Calibration **Sample Name** P2 a cal 6-250ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.716 | 12086700 | 1991265 | 6.0699 | 244.2761 |
| Zopiclone | Zopiclone-D4 | 6.131 | 2383982 | 312362 | 7.6321 | 240.9017 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 449844 | 67916 | 6.6236 | 272.4341 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 267835 | 55331 | 4.8406 | 262.7189 |
| Midazolam | Midazolam-D4 | 7.543 | 2407067 | 1722499 | 1.3974 | 242.8699 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 4040968 | 1283810 | 3.1476 | 251.3577 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 2366292 | 1595949 | 1.4827 | 248.8917 |

15

ISP FORENSICS - Pocatello Instrument # 59740

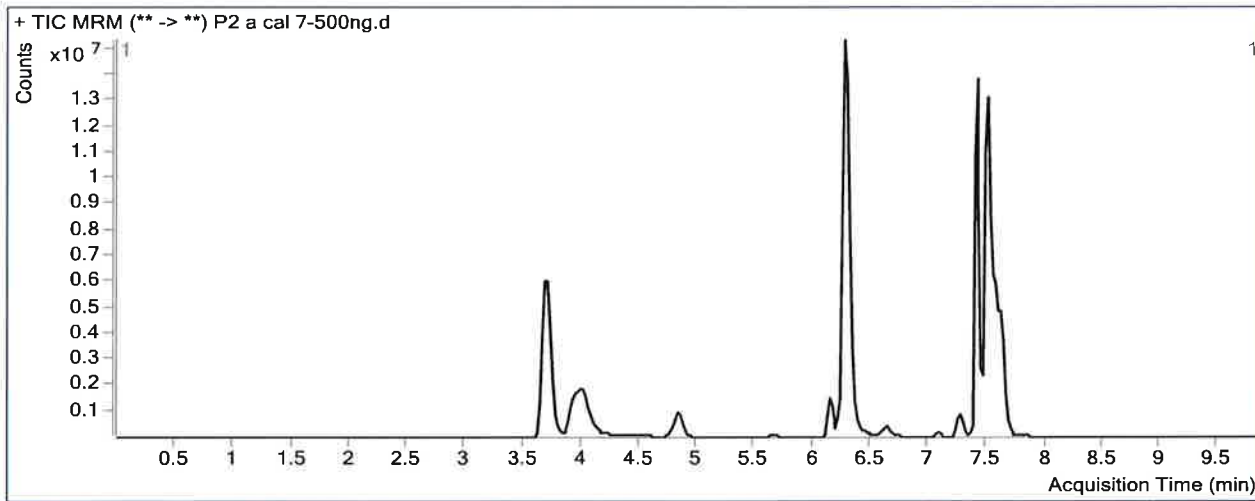
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 18:34 **Data File** P2 a cal 7-500ng.d
Sample Type Calibration **Sample Name** P2 a cal 7-500ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.716 | 24847444 | 1949431 | 12.7460 | 510.9553 |
| Zopiclone | Zopiclone-D4 | 6.131 | 3847825 | 238251 | 16.1503 | 506.9041 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 1261763 | 104650 | 12.0570 | 495.0518 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 705312 | 78095 | 9.0315 | 489.1170 |
| Midazolam | Midazolam-D4 | 7.543 | 4511193 | 1576224 | 2.8620 | 495.3862 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 6444912 | 988550 | 6.5196 | 519.5259 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.605 | 3735944 | 1250760 | 2.9869 | 499.4847 |

B

ISP FORENSICS - Pocatello Instrument # 59740

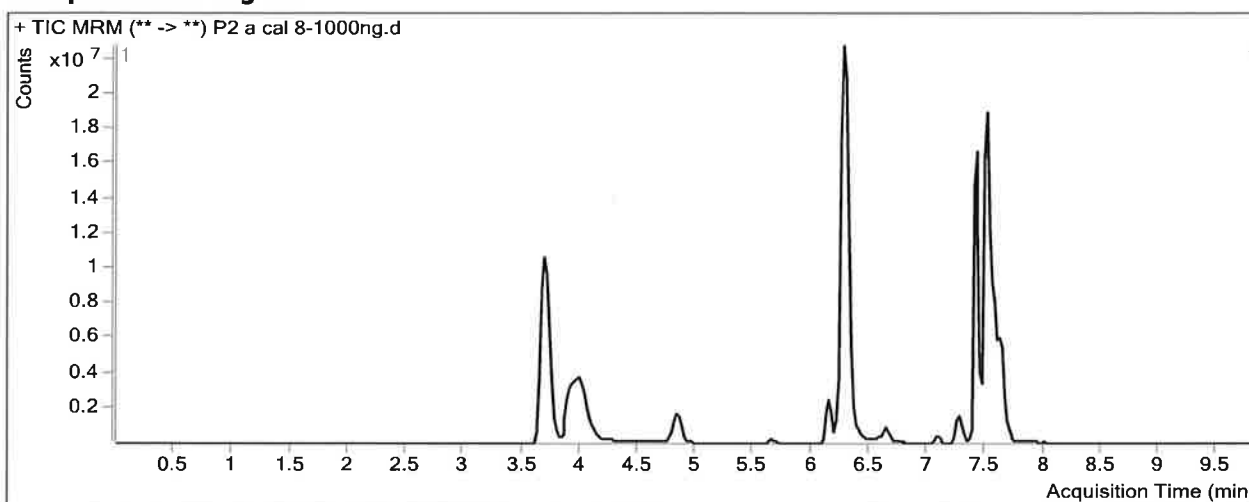
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2a MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:37 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:39 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:37 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 19:00 **Data File** P2 a cal 8-1000ng.d
Sample Type Calibration **Sample Name** P2 a cal 8-1000ng.d
Dilution 1 **Acq Method** MDQ Panel 2a 6-18-18.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 | 3.696 | 46718834 | 1861592 | 25.0962 | 1004.2843 |
| Zopiclone | Zopiclone-D4 | 6.131 | 6247062 | 192934 | 32.3792 | 1013.6964 |
| Amitriptyline | Amitriptyline-D3 | 7.489 | 1975342 | 79898 | 24.7232 | 1014.0148 |
| Nortriptyline | Nortriptyline-D3 | 7.514 | 1124028 | 47160 | 23.8344 | 1288.7861 |
| Midazolam | Midazolam-D4 | 7.543 | 7558562 | 1280468 | 5.9030 | 1019.6848 |
| alpha-hydroxymidazolam | alpha-hydroxymidazolam-D4 | 7.562 | 8694563 | 684778 | 12.6969 | 1010.8092 |
| Chlordiazepoxide | Chlordiazepoxide-D5 | 7.625 | 5289771 | 874384 | 6.0497 | 1009.7105 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

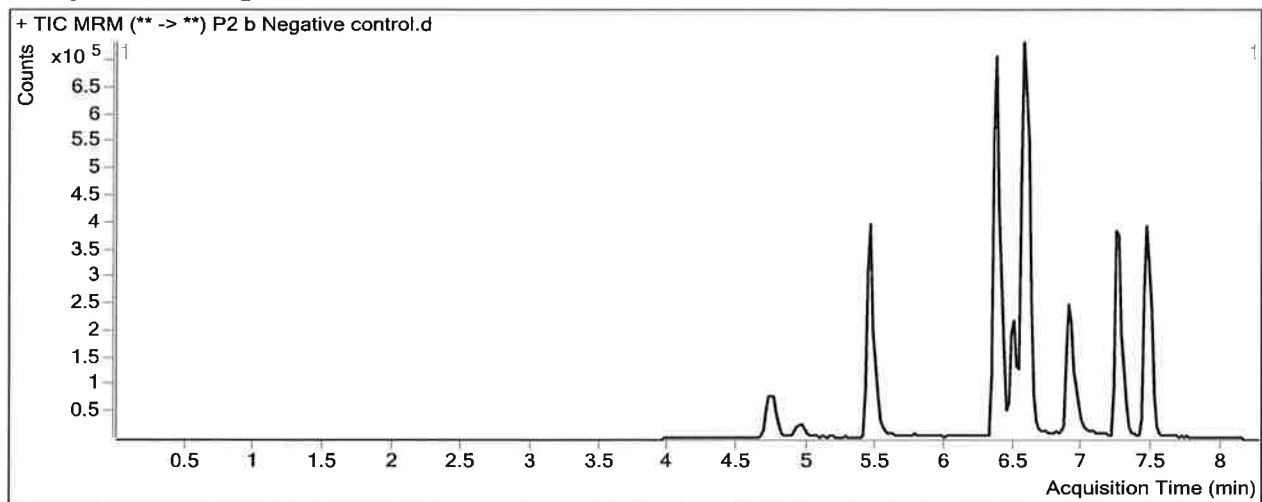
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 fi
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 19:51 **Data File** P2 b Negative control.d
Sample Type Sample **Sample Name** P2 b Negative control.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.m
Position P1-E6 **Sample Info**
Inj Vol -1 **Comment** AM 28-Panel 2

Sample Chromatogram



15

ISP FORENSICS - Pocatello Instrument # 59740

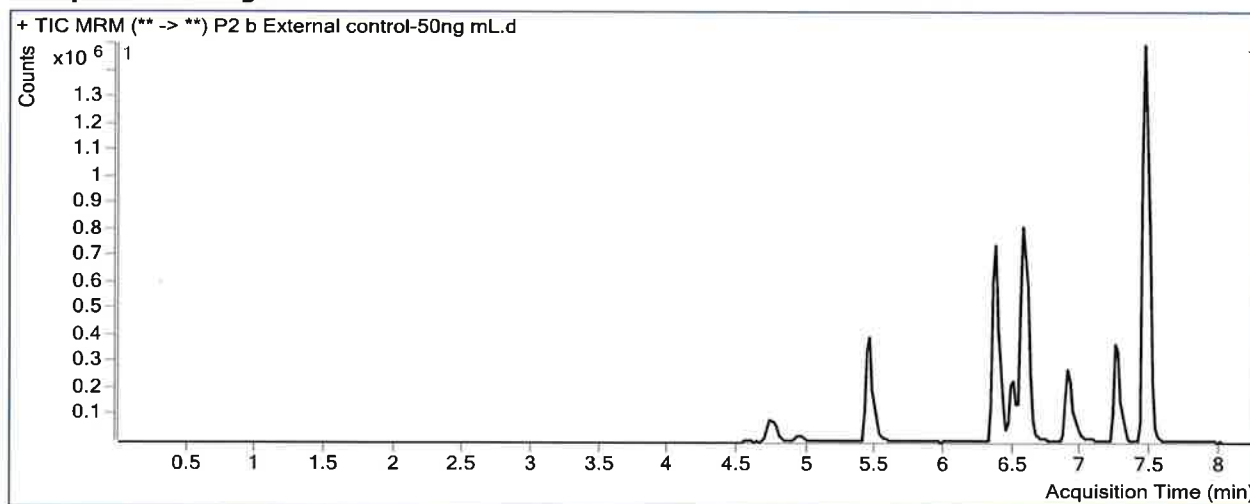
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 20:16 **Data File** P2 b External control-50ng mL.d
Sample Type Sample **Sample Name** P2 b External control-50ng mL.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.m
Position P1-F6 **Sample Info**
Inj Vol -1 **Comment** AM 28; WS061218 50 ng flunitrazepam, zopiclone, chlorodiazepoxide

Sample Chromatogram



Results

| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|----------------------|------------------|-------|----------|-----------|------------|----------------|
| <u>Flunitrazepam</u> | Flunitrazepam-D7 | 7.456 | 2023246 | 556165 | 3.6379 | <u>39.8121</u> |

TS

ISP FORENSICS - Pocatello Instrument # 59740

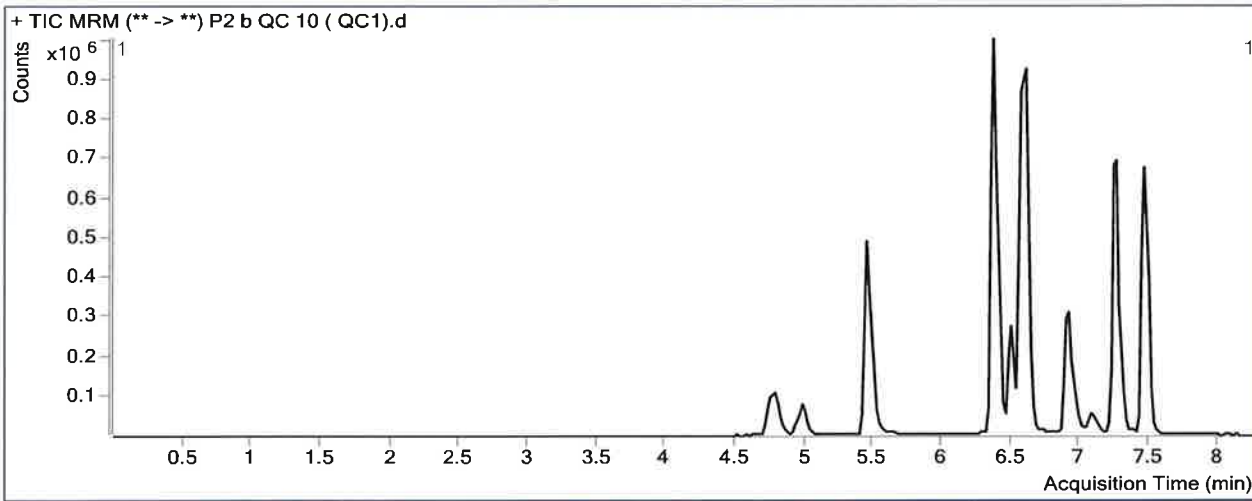
Multi-Drug Confirmatory Analysis Report

| | | | |
|--------------------------|---|----------------------|-----------|
| Batch Data Path | C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir | | |
| Analysis Time | 6/22/2018 1:41 PM | Analyst Name | ISPUser |
| Report Time | 6/22/2018 1:42 PM | Reporter Name | ISPUser |
| Last Calib Update | 6/22/2018 1:41 PM | Batch State | Processed |

Analysis Info

| | | | |
|--------------------|------------------|--------------------|-------------------------|
| Acq Time | 2018-06-18 21:07 | Data File | P2 b QC 10 (QC1).d |
| Sample Type | QC | Sample Name | P2 b QC 10 (QC1).d |
| Dilution | 1 | Acq Method | MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 454030 | 537168 | 0.8452 | 11.1250 |

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ISP FORENSICS - Pocatello Instrument # 59740

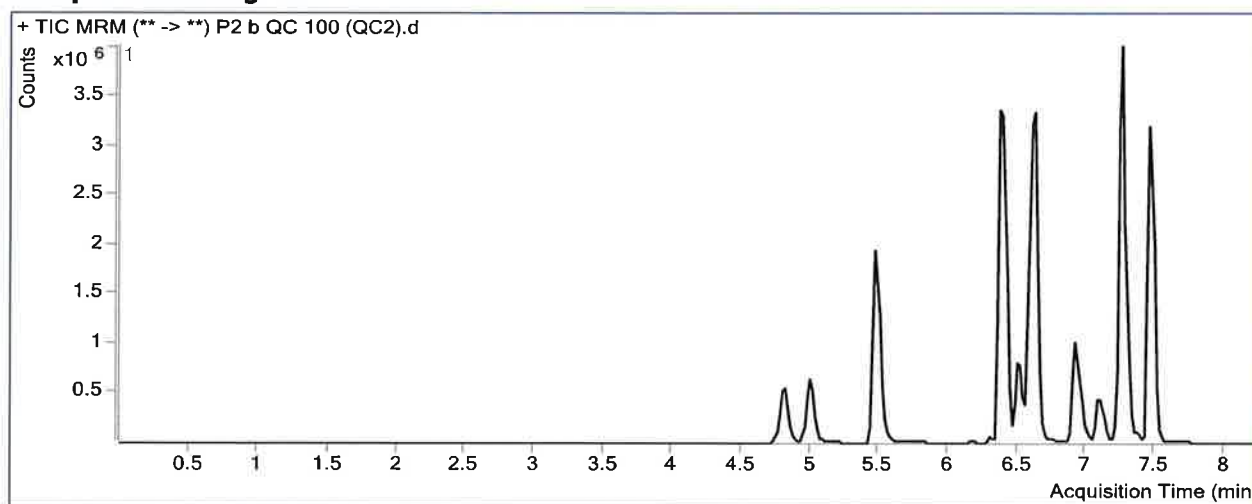
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 22:24 **Data File** P2 b QC 100 (QC2).d
Sample Type QC **Sample Name** P2 b QC 100 (QC2).d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 4323153 | 479032 | 9.0248 | 95.1488 |

B

ISP FORENSICS - Pocatello Instrument # 59740

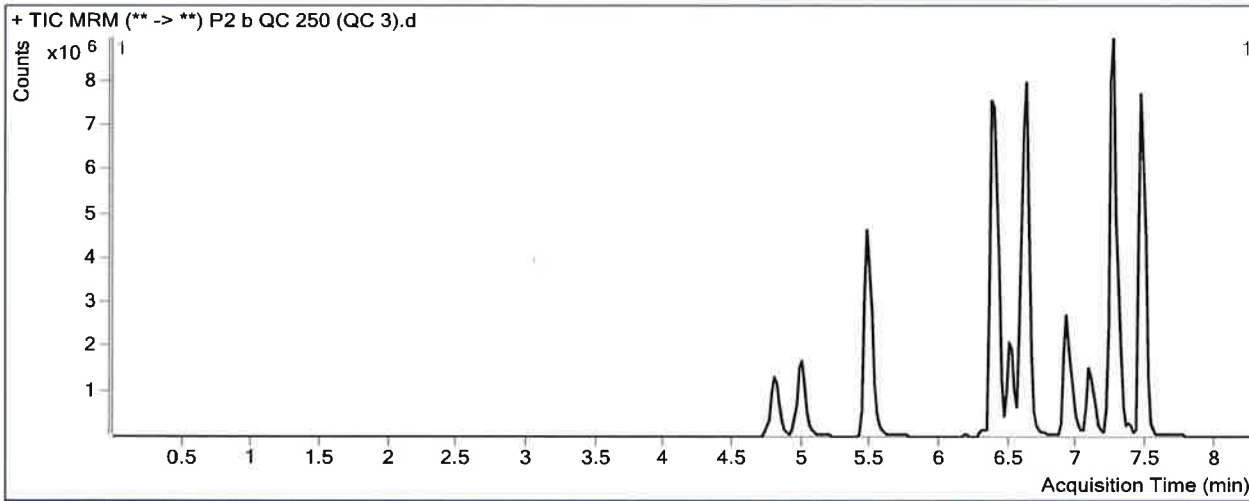
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 23:40 **Data File** P2 b QC 250 (QC 3).d
Sample Type QC **Sample Name** P2 b QC 250 (QC 3).d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 9319710 | 394085 | 23.6490 | 245.3754 |

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ISP FORENSICS - Pocatello Instrument # 59740

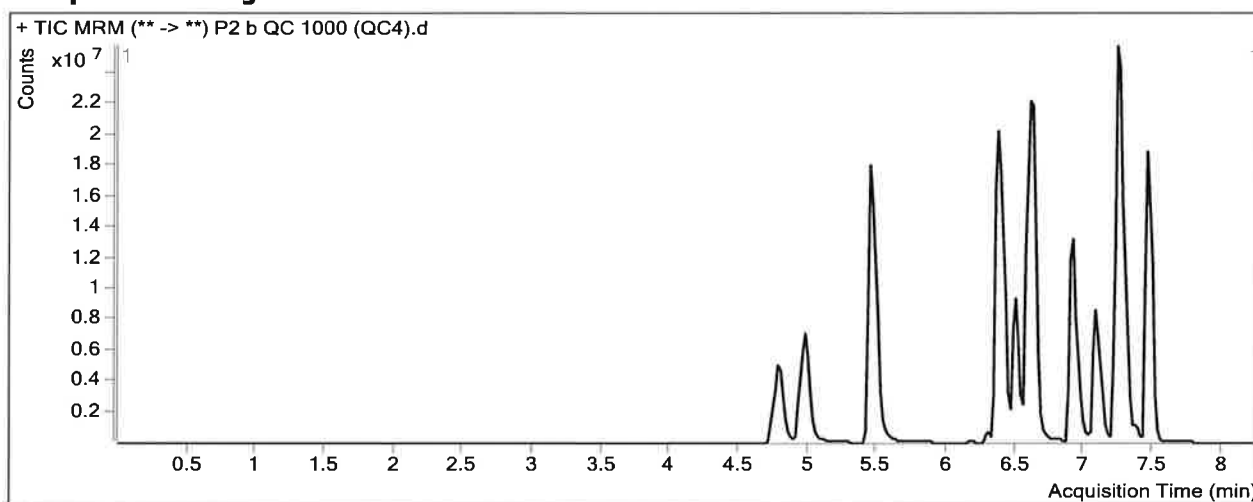
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-19 00:56 **Data File** P2 b QC 1000 (QC4).d
Sample Type QC **Sample Name** P2 b QC 1000 (QC4).d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.m
Position P1-D6 **Sample Info**
Inj Vol -1 **Comment** AM 28-Panel 2

Sample Chromatogram



Results

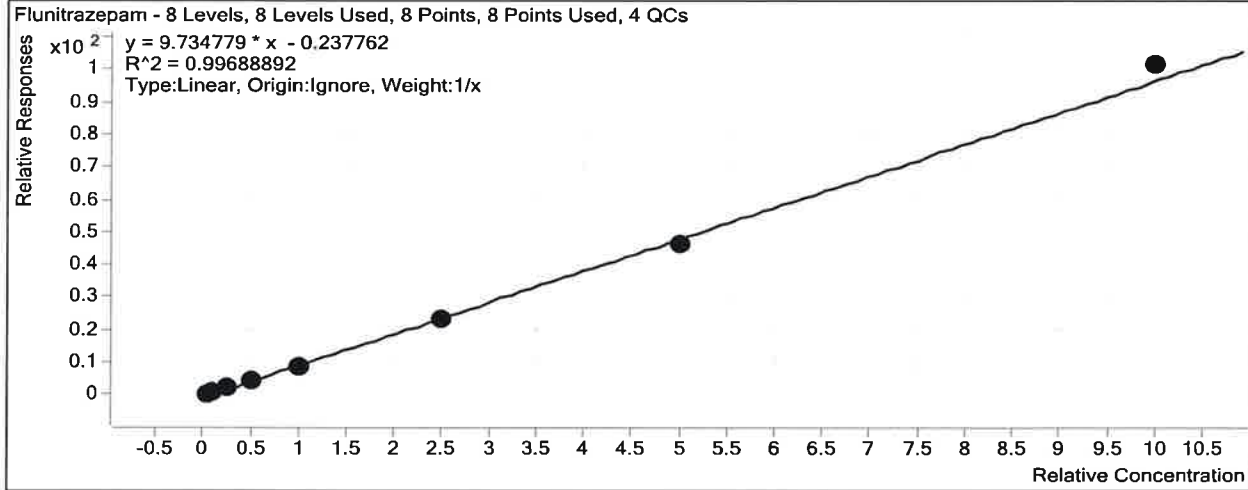
| Compound | ISTD Compound | RT | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 20621214 | 201548 | 102.3144 | 1053.4617 |

ISP Forensics Calibration Curve Report

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Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bin
Last Calib Update 6/22/2018 1:41 PM **Analyst Name** ISP TOX

Target Compound Flunitrazepam
Internal Standard Flunitrazepam-D7



| Sample | Level | Enabled | Exp Conc | Final Conc | Accuracy |
|----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 b cal 1-5ng.d | 1 | <input checked="" type="checkbox"/> | 5 | 6.4 | 127.4 |
| P2 b cal 2-10ng.d | 2 | <input checked="" type="checkbox"/> | 10 | 10.2 | 102.4 |
| P2 b QC 10 (QC1).d | 2 | <input checked="" type="checkbox"/> | 10 | 11.1 | 111.2 |
| P2 b cal 3-25ng.d | 3 | <input checked="" type="checkbox"/> | 25 | 23.9 | 95.4 |
| P2 b cal 4-50ng.d | 4 | <input checked="" type="checkbox"/> | 50 | 44.9 | 89.8 |
| P2 b cal 5-100ng.d | 5 | <input checked="" type="checkbox"/> | 100 | 87.3 | 87.3 |
| P2 b QC 100 (QC2).d | 5 | <input checked="" type="checkbox"/> | 100 | 95.1 | 95.1 |
| P2 b cal 6-250ng.d | 6 | <input checked="" type="checkbox"/> | 250 | 243.1 | 97.2 |
| P2 b QC 250 (QC 3).d | 6 | <input checked="" type="checkbox"/> | 250 | 245.4 | 98.2 |
| P2 b cal 7-500ng.d | 7 | <input checked="" type="checkbox"/> | 500 | 480.0 | 96.0 |
| P2 b cal 8-1000ng.d | 8 | <input checked="" type="checkbox"/> | 1000 | 1044.3 | 104.4 |
| P2 b QC 1000 (QC4).d | 8 | <input checked="" type="checkbox"/> | 1000 | 1053.5 | 105.3 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

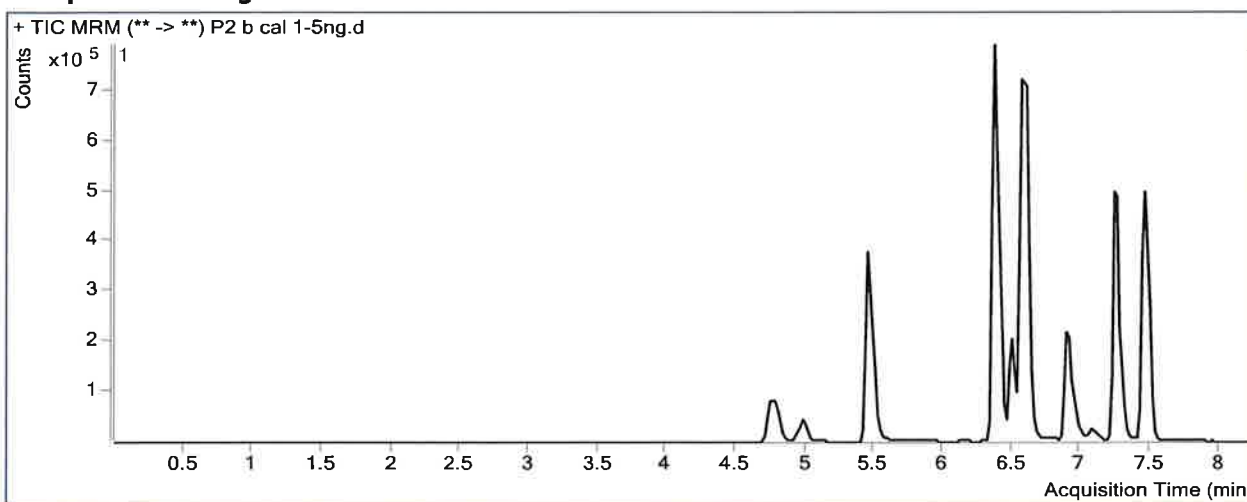
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 16:14 **Data File** P2 b cal 1-5ng.d
Sample Type Calibration **Sample Name** P2 b cal 1-5ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 209930 | 548824 | 0.3825 | 6.3717 |

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ISP FORENSICS - Pocatello Instrument # 59740

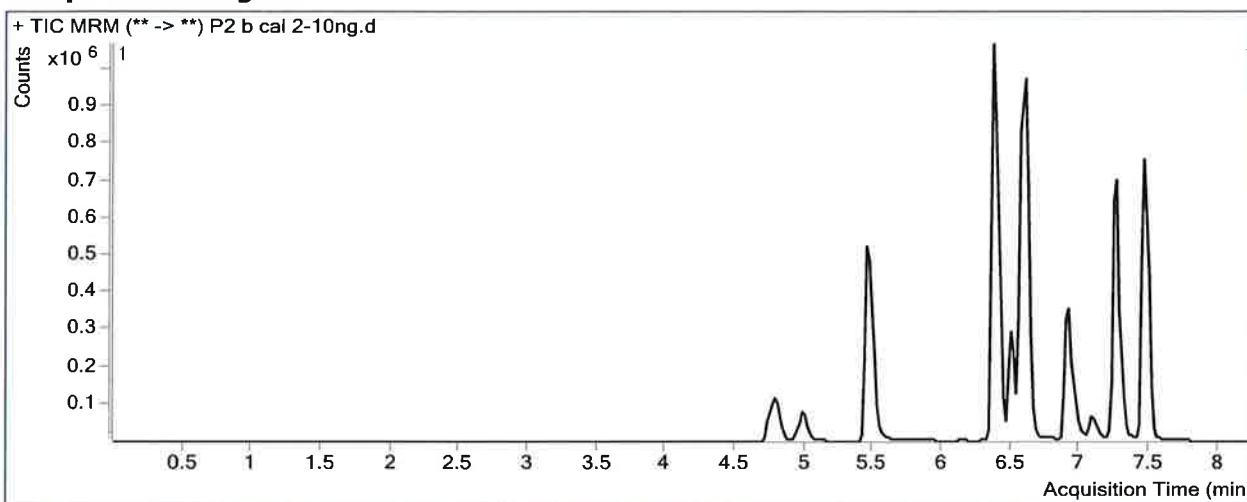
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 16:39 **Data File** P2 b cal 2-10ng.d
Sample Type Calibration **Sample Name** P2 b cal 2-10ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 437512 | 576543 | 0.7589 | 10.2377 |

B

ISP FORENSICS - Pocatello Instrument # 59740

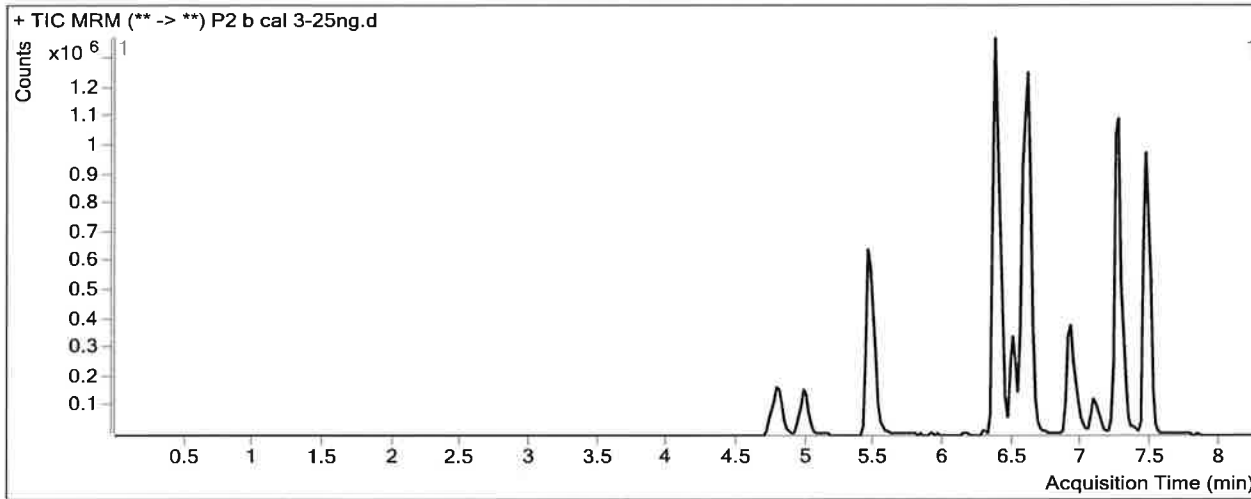
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 17:05 **Data File** P2 b cal 3-25ng.d
Sample Type Calibration **Sample Name** P2 b cal 3-25ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 1033729 | 495770 | 2.0851 | 23.8615 |

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ISP FORENSICS - Pocatello Instrument # 59740

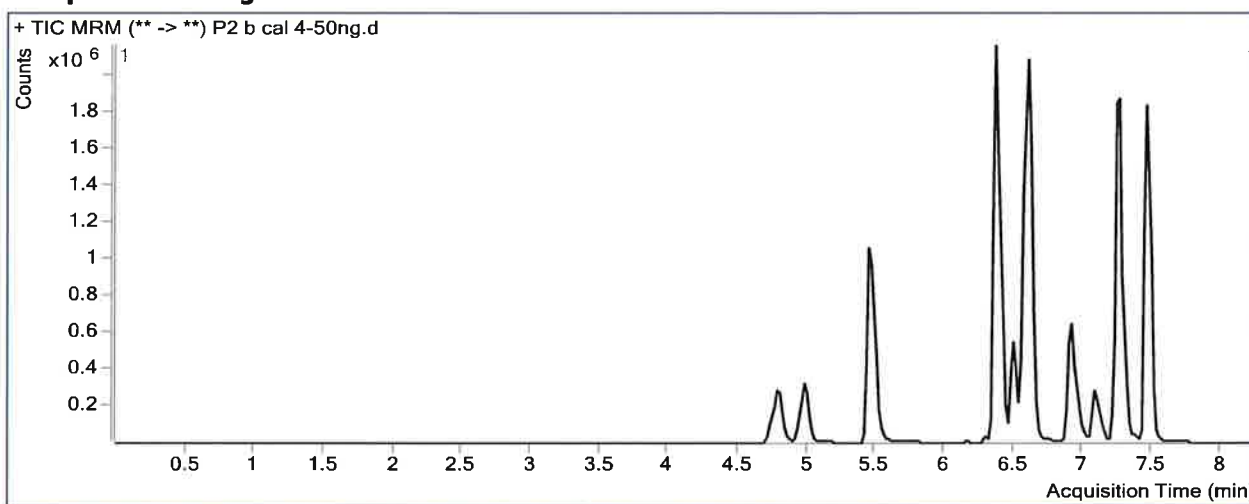
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 17:30 **Data File** P2 b cal 4-50ng.d
Sample Type Calibration **Sample Name** P2 b cal 4-50ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 2041585 | 493860 | 4.1339 | 44.9081 |

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ISP FORENSICS - Pocatello Instrument # 59740

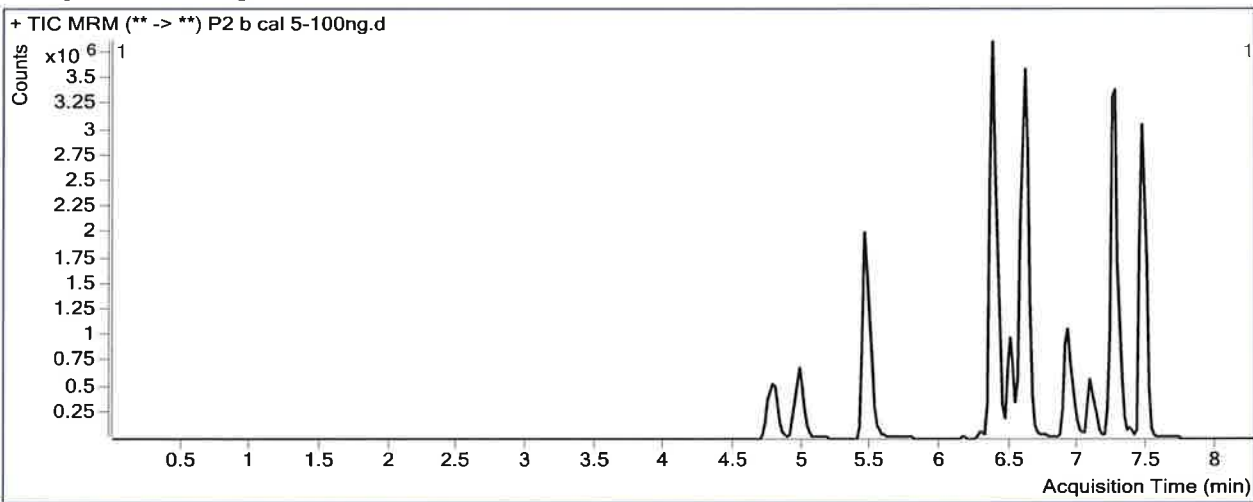
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 17:56 **Data File** P2 b cal 5-100ng.d
Sample Type Calibration **Sample Name** P2 b cal 5-100ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 3820115 | 462578 | 8.2583 | 87.2755 |

TS

ISP FORENSICS - Pocatello Instrument # 59740

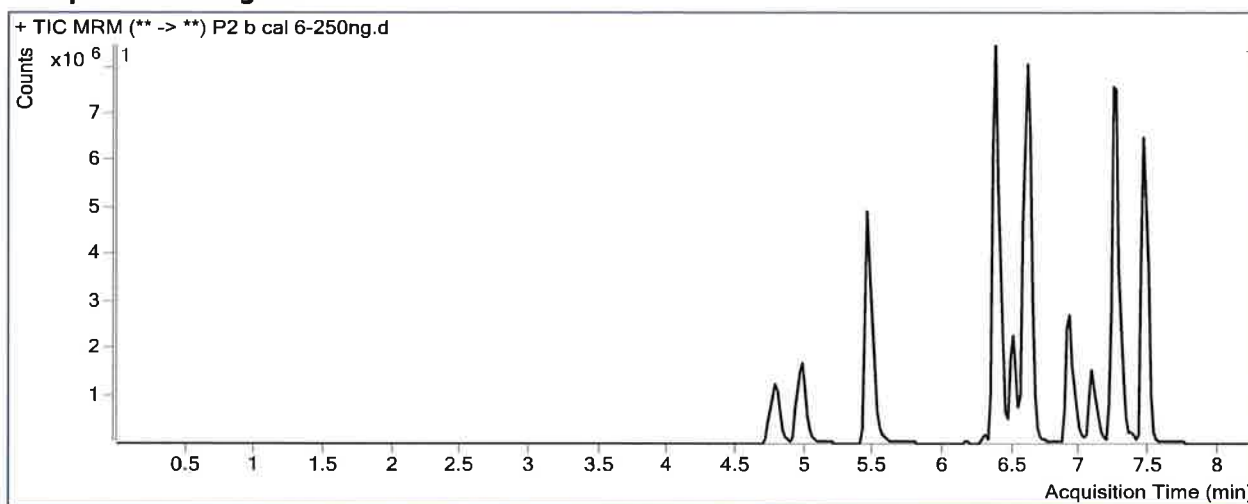
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 18:21 **Data File** P2 b cal 6-250ng.d
Sample Type Calibration **Sample Name** P2 b cal 6-250ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 8470425 | 361629 | 23.4230 | 243.0537 |

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ISP FORENSICS - Pocatello Instrument # 59740

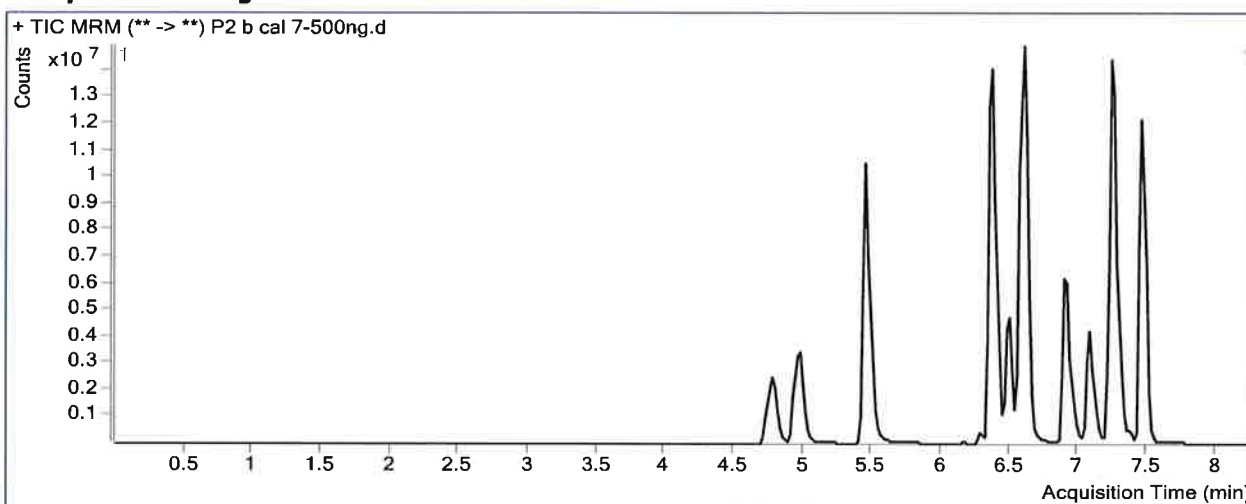
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:41 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 18:47 **Data File** P2 b cal 7-500ng.d
Sample Type Calibration **Sample Name** P2 b cal 7-500ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 13617776 | 292917 | 46.4902 | 480.0102 |

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ISP FORENSICS - Pocatello Instrument # 59740

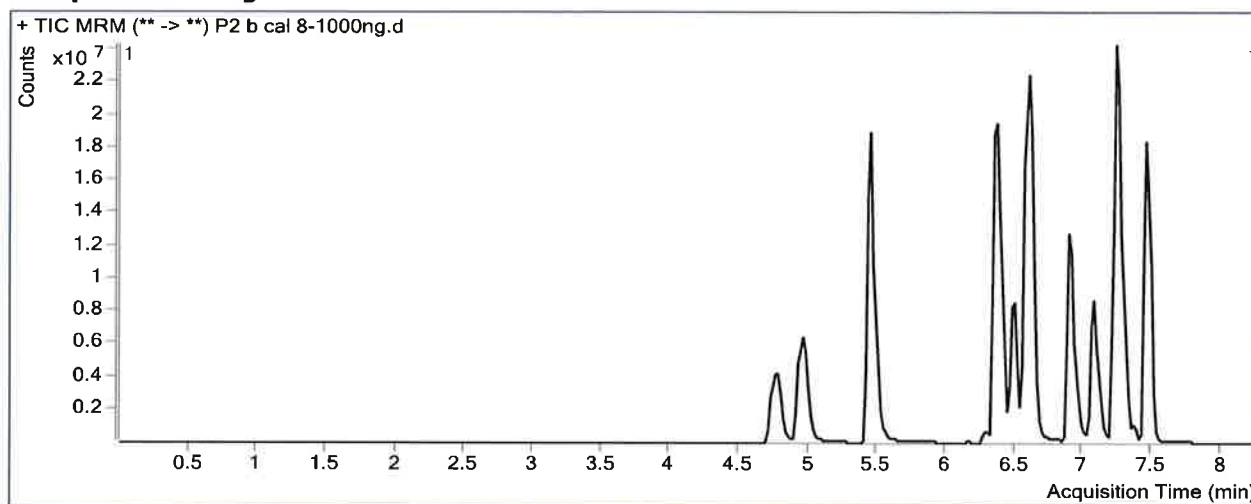
Multi-Drug Confirmatory Analysis Report

Batch Data Path C:\MassHunter\Data\2018\MD Quant\061818 MDQ P1 and P2 TS\QuantResults\p2b MDQ 61818 final.batch.bir
Analysis Time 6/22/2018 1:41 PM **Analyst Name** ISPUser
Report Time 6/22/2018 1:42 PM **Reporter Name** ISPUser
Last Calib Update 6/22/2018 1:41 PM **Batch State** Processed

Analysis Info

Acq Time 2018-06-18 19:12 **Data File** P2 b cal 8-1000ng.d
Sample Type Calibration **Sample Name** P2 b cal 8-1000ng.d
Dilution 1 **Acq Method** MDQ Panel 2b 06-18-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.456 | 19475088 | 192023 | 101.4207 | 1044.2817 |