



**Worklist: 2773**

| <u>LAB_CASE</u> | <u>ITEM</u> | <u>TASK_ID</u> | <u>DESCRIPTION</u>                           |   |
|-----------------|-------------|----------------|--|---|
| P2018-2828      | 3           | 130925         | AM 28 Blood Multi-Drug Quant Panel 2 by LC-Q |  |
| P2018-2932      | 1           | 130926         | AM 28 Blood Multi-Drug Quant Panel 2 by LC-Q |  |

P2018-2932-1 was extracted but not injected. Only P2b data was analyzed. Please note that daylight savings took place during this run. At 2:00 am on 11/4/18, the clock on the computer reset back one hour.

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

Extraction Date: 11/02/2018

Analyst: Celena Shrum

Plate lot#: 0530563

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:** 445283-1

**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: 16** 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\110218 MDQ P1 P2 CS*
- Batch Name: *MDQ P2b 110218 CS*
- 2. Make necessary changes to integration limits
- 3. Integration linear and R<sup>2</sup> 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.

### COMMENTS:

*The only compounds evaluated were Flunitrazepam and Phencyclidine.*



# Idaho State Police Forensic Services

CS

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

**Analyst:** Celena Shrum  
**Extraction Date:** 11/02/2018  
**Worklist Number:** 2773

| <i>Reagent</i>   | <i>Lot Number</i> | <i>Expiration Date</i>             | <i>Date in Service</i>             | <i>Date Out of Service</i> | <i>Initials</i> |
|--|-------------------|------------------------------------|------------------------------------|----------------------------|-----------------|
| ToxBox Benzo/Opiate Quant Plate                                    | 0530563           | 05/21/19                           | 06/05/18                           |                            |                 |
| Negative Blood   | 445283-1          |                                    | 10/25/18                           |                            |                 |
| Methanol External Control Solution                                 | 022018            | 02/20/19                           | 02/20/18                           |                            |                 |
| Blood External Control Solution                                    | 081518            | <del>02/15/19</del><br>02/20/19 CS | 08/15/18                           |                            |                 |
| 0.5M Ammonium Hydroxide  | 082818            | 02/28/19                           | 08/28/18                           |                            |                 |
| Formic Acid (LCMS Grade)   | 095180B           |                                    | 11/15/17                           |                            |                 |
| Water (LCMS Grade)   | 182702            |                                    | 06/21/18                           |                            |                 |
| 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                                    | 082818            |                                    | 08/28/18                           |                            |                 |
| 5mM Ammonium Formate + 0.01% Formic Acid in Water (Mobile Phase A) | 110218            |                                    | <del>11/08/18</del><br>11/02/18 CS |                            |                 |
| 0.01% Formic Acid in Methanol (Mobile Phase B)                     | 100918            |                                    | 10/09/18                           |                            |                 |
| Needle Rinse--75% LCMS MeOH in LCMS Water                          | 100918            |                                    | 10/09/18                           |                            |                 |

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

| <i>Component</i> | <i>Source</i>  | <i>Source Lot Number</i> | <i>Expiration Date</i> |
|------------------|----------------|--------------------------|------------------------|
| 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       |                          |                        |

CS

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

**Blood External Control Solution (Lot: WS081518)**

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 |                        | 022018                   |
| Prepared:                          | 08/15/18               |                          |
| Prepared by:                       | Sarah Pickle           |                          |
| Expires:                           | -02/15/18- 02/20/19 CS |                          |

**0.5M Ammonium Hydroxide (Lot: 082818)**

| <i>Component</i>                         | <i>Source</i>  | <i>Source Lot Number</i> |
|--|----------------|--------------------------|
| Ammonium Hydroxide (ACS grade or better) | Fisher         | 053200                   |
| DI Water                                 | -              | -                        |
| Prepared:                                | 08/28/18       |                          |
| Prepared By:                             | Tamara Salazar |                          |
| Expires:                                 | 02/28/18       |                          |

**20% LCMS Methanol in LCMS Water (Lot: 082818)**

| <i>Component</i>   | <i>Source</i>  | <i>Source Lot Number</i> |
|--------------------|----------------|--------------------------|
| MeOH (LCMS Grade)  | Fisher         | 177145                   |
| Water (LCMS Grade) | Fisher         | 182702                   |
| Prepared:          | 08/28/18       |                          |
| Prepared By:       | Tamara Salazar |                          |

**5mM Ammonium Formate + 0.01% Formic Acid in Water (Lot: ~~100918~~ 11/02/18<sup>CS</sup>)**

| <i>Component</i> | <i>Source</i> | <i>Source Lot Number</i> |
|------------------|---------------|--------------------------|
| Ammonium Formate | Arcos Organic | A0287238                 |
| LCMS Water       | Fisher        | 182702                   |
| Formic Acid      | Fisher        | 095180B                  |
| Prepared:        | 11/02/18      |                          |
| Prepared By:     | Celena Shrum  |                          |

**0.01% Formic Acid in LCMS Methanol (Mobile Phase B) (Lot: 100918)**

| <i>Component</i>  | <i>Source</i>  | <i>Source Lot Number</i> |
|-------------------|----------------|--------------------------|
| Formic Acid       | Fisher         | 095180B                  |
| MeOH (LCMS Grade) | Fisher         | 177145                   |
| Prepared:         | 10/09/18       |                          |
| Prepared By:      | Tamara Salazar |                          |

**Needle Rinse (75% LCMS MeOH in LCMS Water) (Lot: 100918)**

| <i>Component</i>   | <i>Source</i> | <i>Source Lot Number</i> |
|--------------------|---------------|--------------------------|
| MeOH (LCMS Grade)  | Fisher        | 177145                   |
| Water (LCMS Grade) | Fisher        | 182702                   |
| Prepared:          | 10/09/18      |                          |
| Prepared By:       | Sarah Pickle  |                          |

# ISP FORENSICS - Pocatello Instrument # 59740

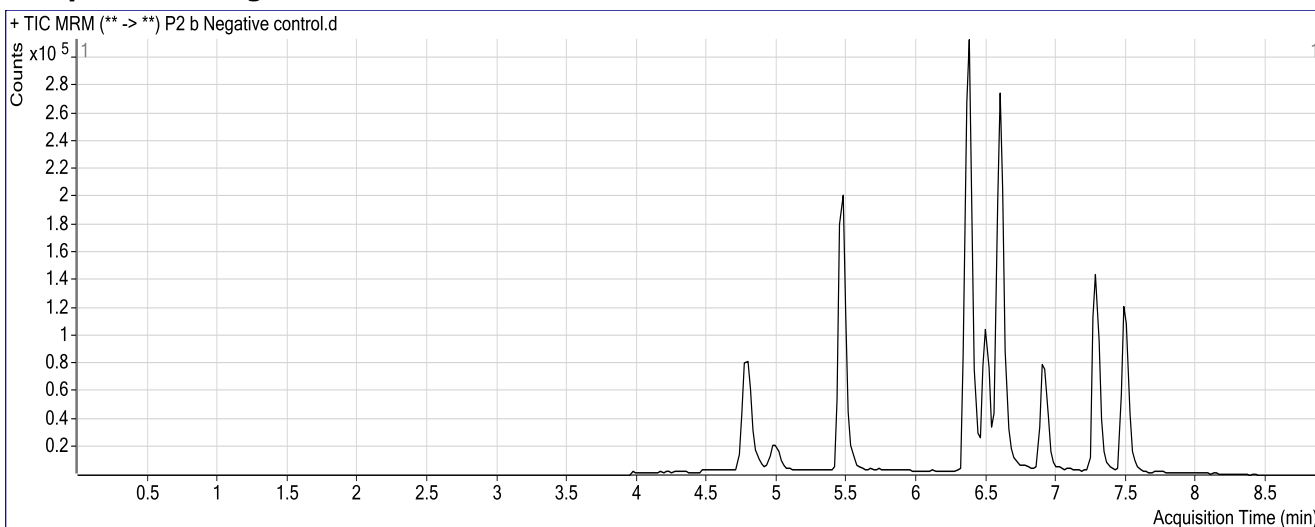
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.t |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

### Analysis Info

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 22:09 | <b>Data File</b>   | P2 b Negative control.d |
| <b>Sample Type</b> | Sample           | <b>Sample Name</b> | P2 b Negative control.d |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-E6            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

### Sample Chromatogram



# ISP FORENSICS - Pocatello Instrument # 59740

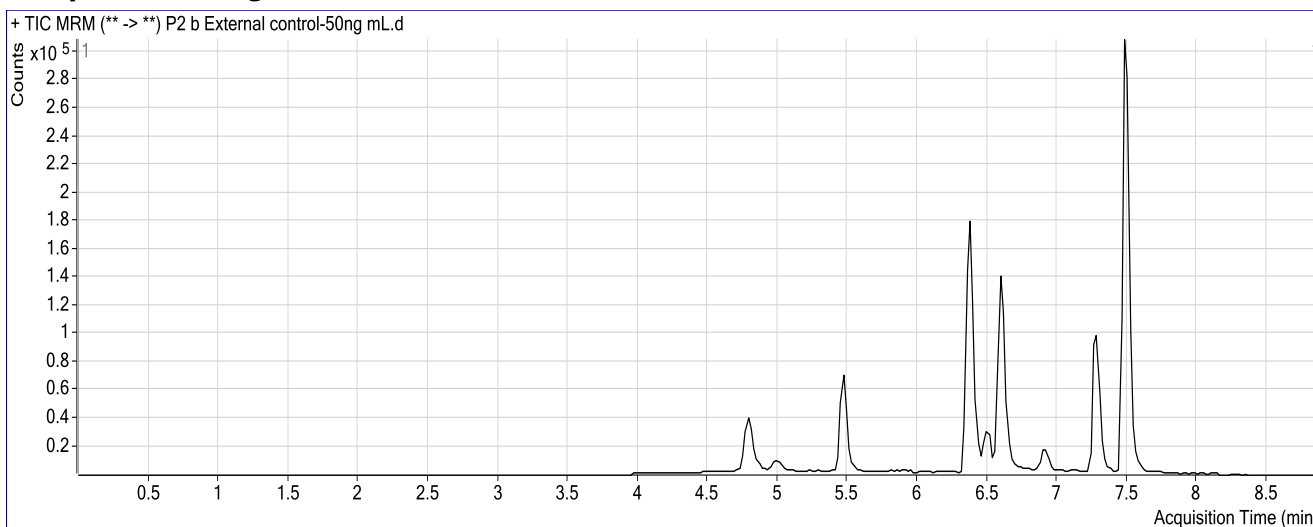
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

### Analysis Info

|                    |                  |                    |   |
|--------------------|------------------|--------------------|---|
| <b>Acq Time</b>    | 2018-11-03 22:35 | <b>Data File</b>   | P2 b External control-50ng mL.d                                       |
| <b>Sample Type</b> | Sample           | <b>Sample Name</b> | P2 b External control-50ng mL.d                                       |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m   |
| <b>Position</b>    | P1-F6            | <b>Sample Info</b> |   |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28; 361331-1 + WS022018 50 ng flunitrazepam, zopiclone, chlorodia. |

### Sample Chromatogram



### Results

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Flunitrazepam | Flunitrazepam-D7 | 7.477 | 582640   | 121162    | 4.8088     | 54.1544    |

# ISP FORENSICS - Pocatello Instrument # 59740

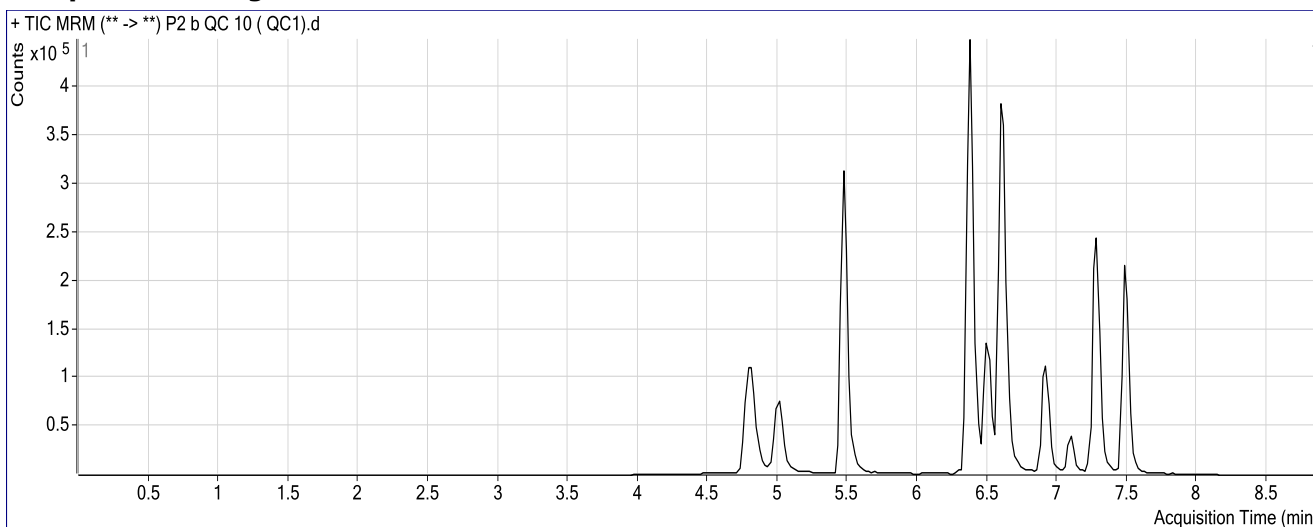
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin  
**Analysis Time** 11/8/2018 2:09 PM **Analyst Name** ISPUser  
**Report Time** 11/9/2018 12:36 PM **Reporter Name** ISPUser  
**Last Calib Update** 11/8/2018 2:09 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-11-03 23:26 **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 06-22-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 70855    | 318794    | 0.2223     | 9.3827     |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 124424   | 170453    | 0.7300     | 9.0595     |

# ISP FORENSICS - Pocatello Instrument # 59740

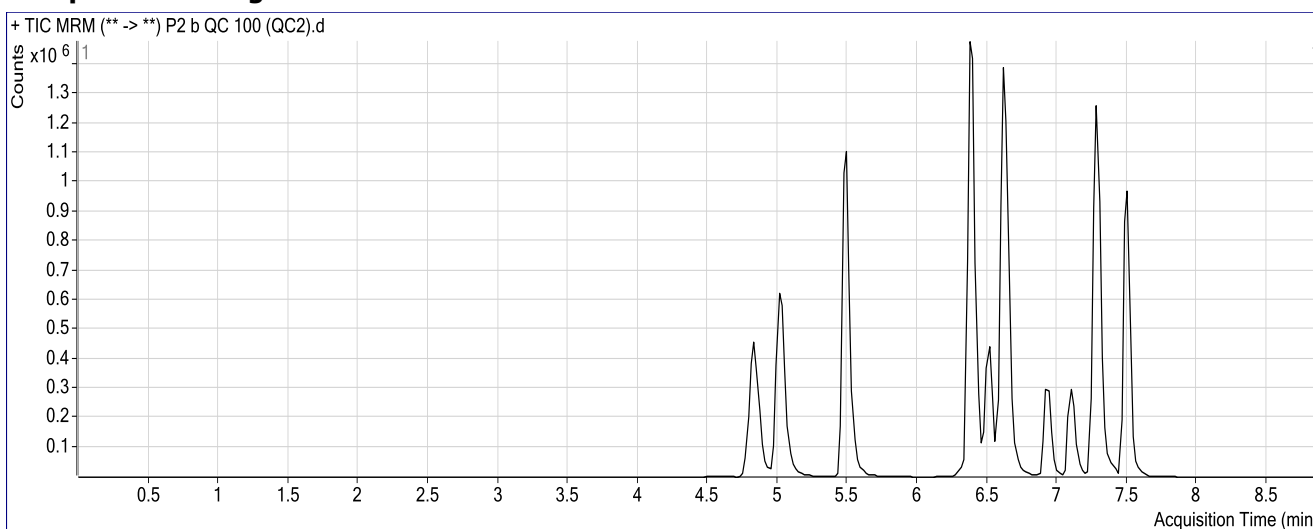
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin  
**Analysis Time** 11/8/2018 2:09 PM **Analyst Name** ISPUser  
**Report Time** 11/9/2018 12:36 PM **Reporter Name** ISPUser  
**Last Calib Update** 11/8/2018 2:09 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-11-04 00:17 **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 06-22-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.947 | 613846   | 236069    | 2.6003     | 97.3720    |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 1383700  | 153598    | 9.0086     | 100.5869   |



# ISP FORENSICS - Pocatello Instrument # 59740

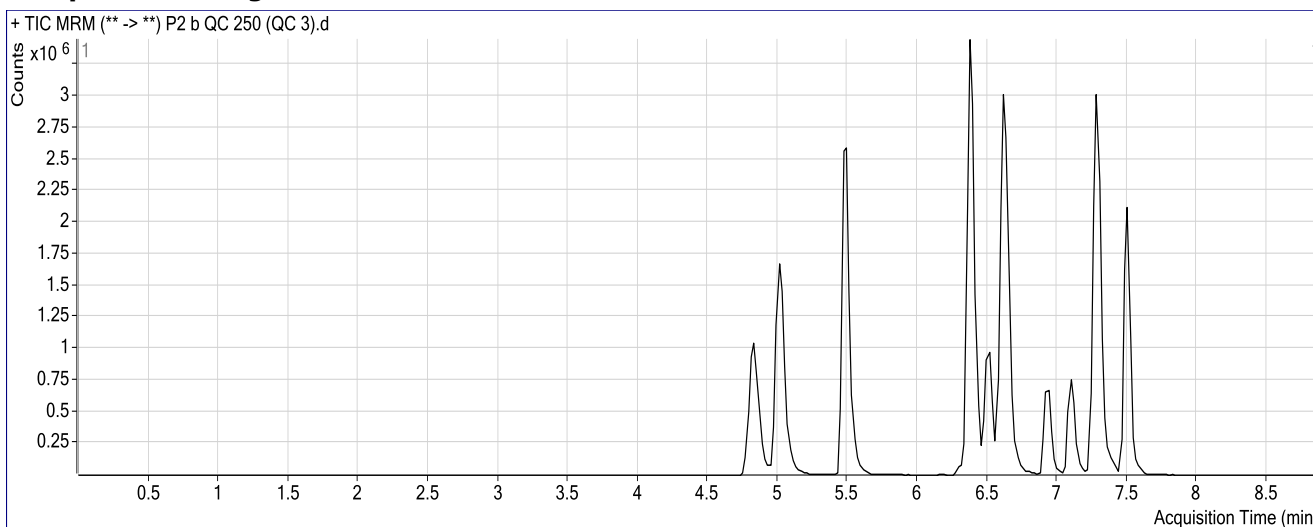
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:37 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

### Analysis Info

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-04 01:33 | <b>Data File</b>   | P2 b QC 250 (QC 3).d    |
| <b>Sample Type</b> | QC               | <b>Sample Name</b> | P2 b QC 250 (QC 3).d    |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-C6            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

### Sample Chromatogram



### Results

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.947 | 1579333  | 233844    | 6.7538     | 251.0552   |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 3210403  | 131274    | 24.4557    | 271.3687   |

# ISP FORENSICS - Pocatello Instrument # 59740

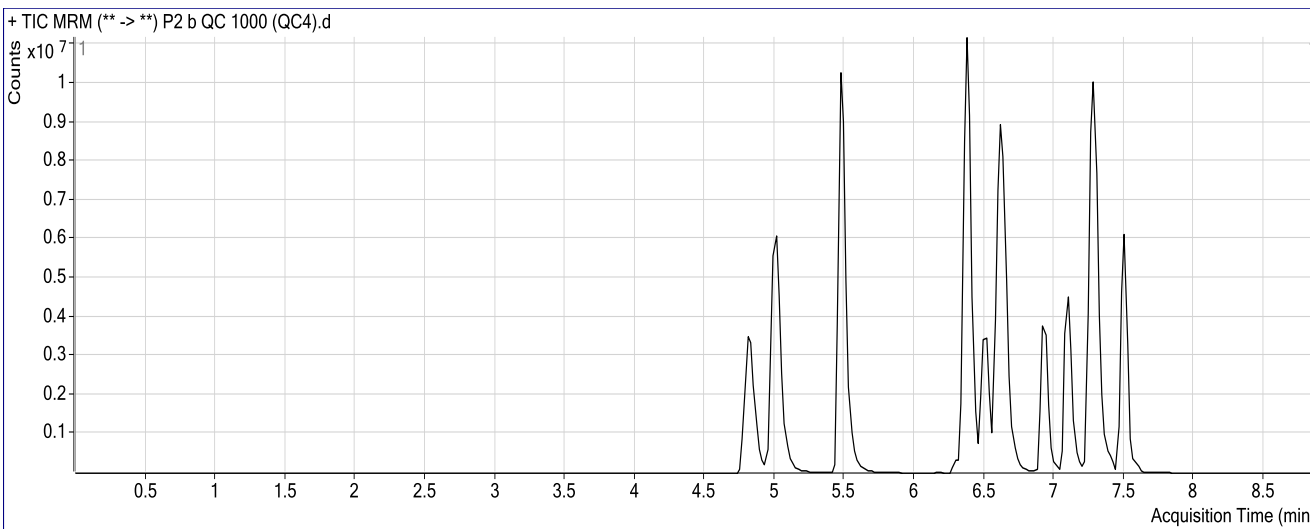
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:37 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-04 01:49 | <b>Data File</b>   | P2 b QC 1000 (QC4).d    |
| <b>Sample Type</b> | QC               | <b>Sample Name</b> | P2 b QC 1000 (QC4).d    |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-D6            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

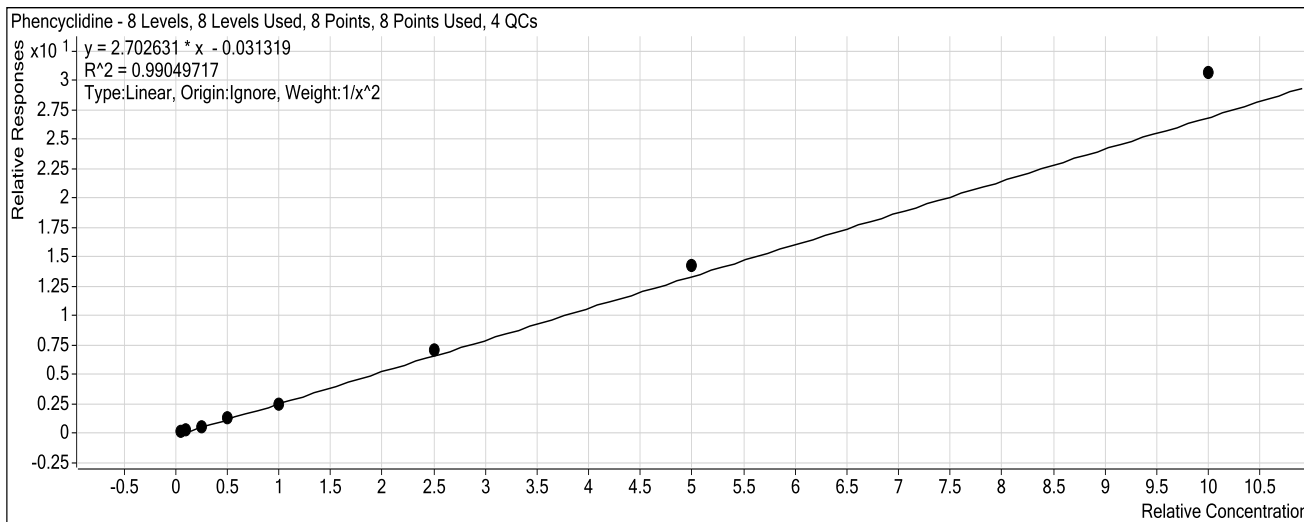
| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 9178154  | 298889    | 30.7076    | 1137.3707  |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 6787873  | 65704     | 103.3103   | 1143.1756  |

# ISP Forensics Calibration Curve Report

CS

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin  
**Last Calib Update** 11/8/2018 2:09 PM **Analyst Name** ISP TOX

**Target Compound** Phencyclidine  
**Internal Standard** Phencyclidine-D5



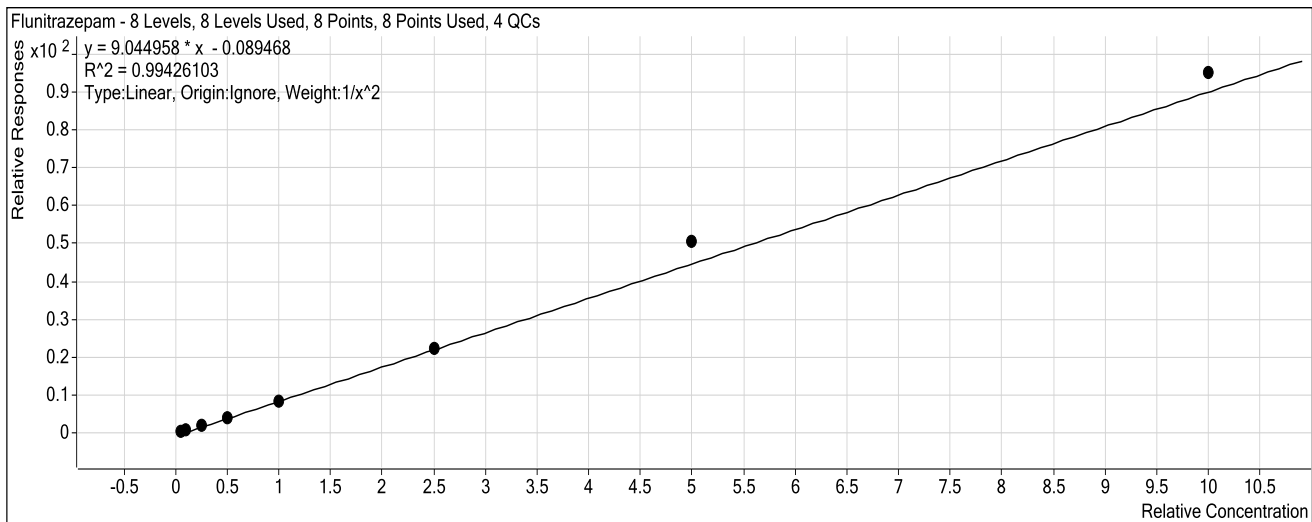
| Sample               | Level | Enabled                             | Exp Conc | Final Conc | Accuracy |
|----------------------|-------|-------------------------------------|----------|------------|----------|
| P2 b cal 1-5ng.d     | 1     | <input checked="" type="checkbox"/> | 5        | 5.3        | 105.7    |
| P2 b cal 2-10ng.d    | 2     | <input checked="" type="checkbox"/> | 10       | 9.4        | 94.3     |
| P2 b QC 10 ( QC1).d  | 2     | <input checked="" type="checkbox"/> | 10       | 9.4        | 93.8     |
| P2 b cal 3-25ng.d    | 3     | <input checked="" type="checkbox"/> | 25       | 22.5       | 89.9     |
| P2 b cal 4-50ng.d    | 4     | <input checked="" type="checkbox"/> | 50       | 46.5       | 93.1     |
| P2 b cal 5-100ng.d   | 5     | <input checked="" type="checkbox"/> | 100      | 93.0       | 93.0     |
| P2 b QC 100 (QC2).d  | 5     | <input checked="" type="checkbox"/> | 100      | 97.4       | 97.4     |
| P2 b cal 6-250ng.d   | 6     | <input checked="" type="checkbox"/> | 250      | 261.2      | 104.5    |
| P2 b QC 250 (QC 3).d | 6     | <input checked="" type="checkbox"/> | 250      | 251.1      | 100.4    |
| P2 b cal 7-500ng.d   | 7     | <input checked="" type="checkbox"/> | 500      | 529.1      | 105.8    |
| P2 b cal 8-1000ng.d  | 8     | <input checked="" type="checkbox"/> | 1000     | 1137.4     | 113.7    |
| P2 b QC 1000 (QC4).d | 8     | <input checked="" type="checkbox"/> | 1000     | 1137.4     | 113.7    |

# ISP Forensics Calibration Curve Report

CS

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin  
**Last Calib Update** 11/8/2018 2:09 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        | 5.2        | 103.3    |
| P2 b cal 2-10ng.d    | 2     | <input checked="" type="checkbox"/> | 10       | 9.7        | 97.4     |
| P2 b QC 10 ( QC1).d  | 2     | <input checked="" type="checkbox"/> | 10       | 9.1        | 90.6     |
| P2 b cal 3-25ng.d    | 3     | <input checked="" type="checkbox"/> | 25       | 23.5       | 94.2     |
| P2 b cal 4-50ng.d    | 4     | <input checked="" type="checkbox"/> | 50       | 46.5       | 93.0     |
| P2 b cal 5-100ng.d   | 5     | <input checked="" type="checkbox"/> | 100      | 95.3       | 95.3     |
| P2 b QC 100 (QC2).d  | 5     | <input checked="" type="checkbox"/> | 100      | 100.6      | 100.6    |
| P2 b cal 6-250ng.d   | 6     | <input checked="" type="checkbox"/> | 250      | 247.9      | 99.2     |
| P2 b QC 250 (QC 3).d | 6     | <input checked="" type="checkbox"/> | 250      | 271.4      | 108.5    |
| P2 b cal 7-500ng.d   | 7     | <input checked="" type="checkbox"/> | 500      | 561.7      | 112.3    |
| P2 b cal 8-1000ng.d  | 8     | <input checked="" type="checkbox"/> | 1000     | 1053.7     | 105.4    |
| P2 b QC 1000 (QC4).d | 8     | <input checked="" type="checkbox"/> | 1000     | 1143.2     | 114.3    |

# ISP FORENSICS - Pocatello Instrument # 59740

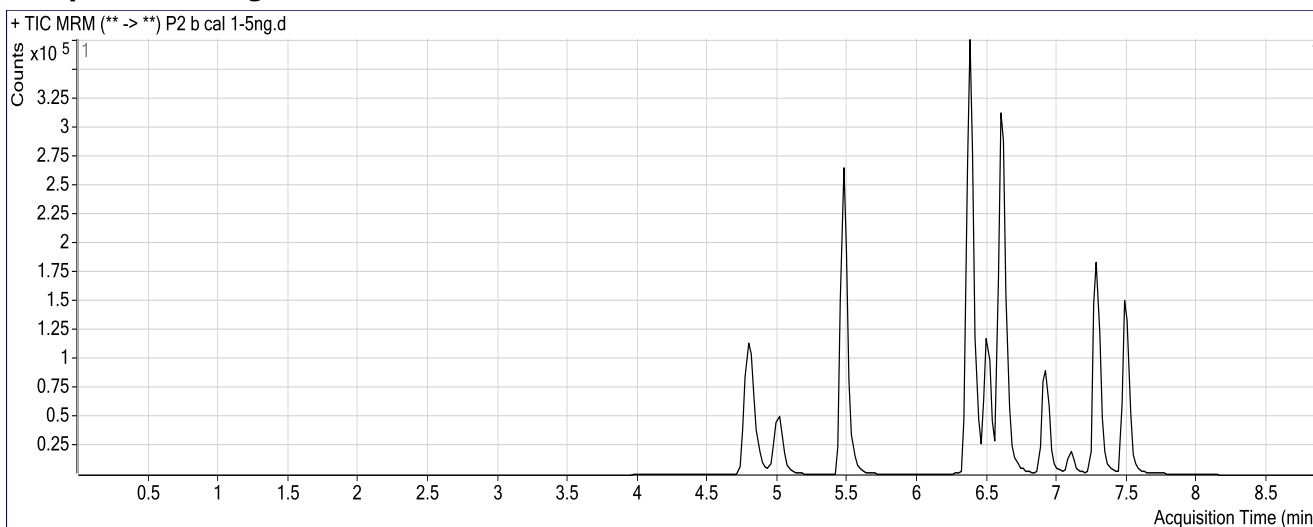
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:35 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

### Analysis Info

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 18:19 | <b>Data File</b>   | P2 b cal 1-5ng.d        |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 1-5ng.d        |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-A5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

### Sample Chromatogram



### Results

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 32951    | 295590    | 0.1115     | 5.2835     |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 58982    | 156238    | 0.3775     | 5.1629     |

# ISP FORENSICS - Pocatello Instrument # 59740

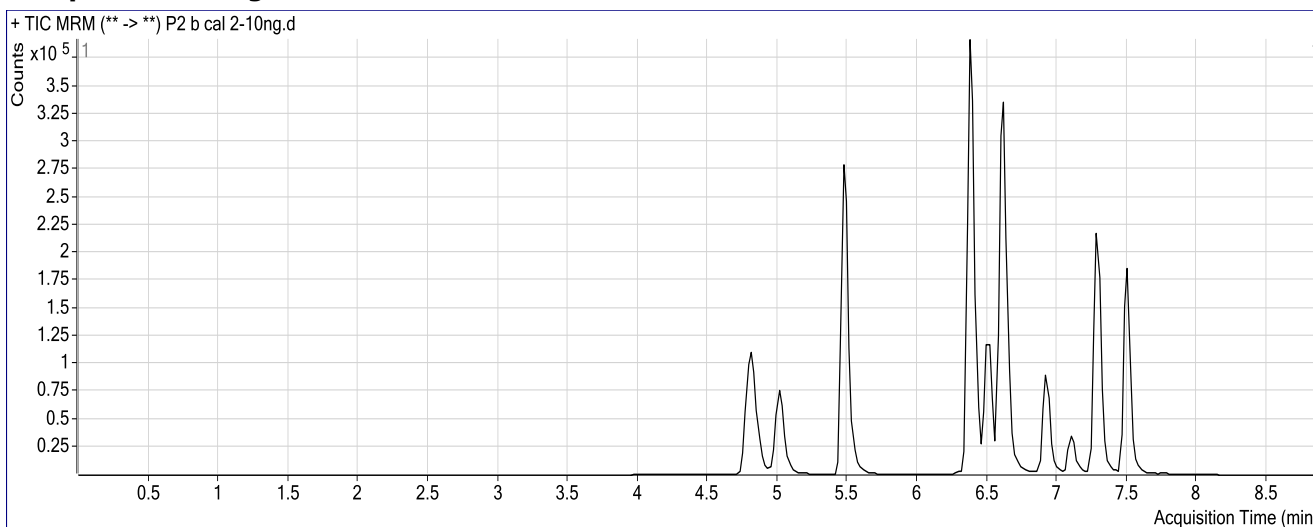
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin  
**Analysis Time** 11/8/2018 2:09 PM **Analyst Name** ISPUser  
**Report Time** 11/9/2018 12:35 PM **Reporter Name** ISPUser  
**Last Calib Update** 11/8/2018 2:09 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-11-03 18:45 **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 06-22-18.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 |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.947 | 54172    | 242197    | 0.2237     | 9.4348     |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 113445   | 143290    | 0.7917     | 9.7423     |

# ISP FORENSICS - Pocatello Instrument # 59740

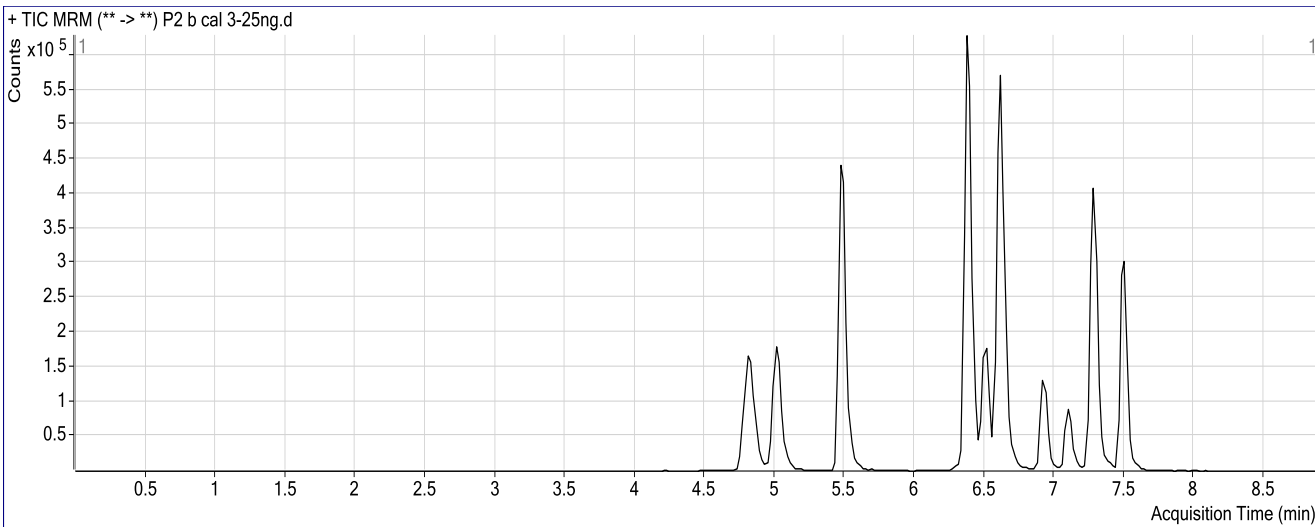
## Multi-Drug Confirmatory Analysis Report

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 19:10 | <b>Data File</b>   | P2 b cal 3-25ng.d       |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 3-25ng.d       |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-C5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.947 | 152319   | 264307    | 0.5763     | 22.4823    |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 336612   | 164994    | 2.0402     | 23.5448    |

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

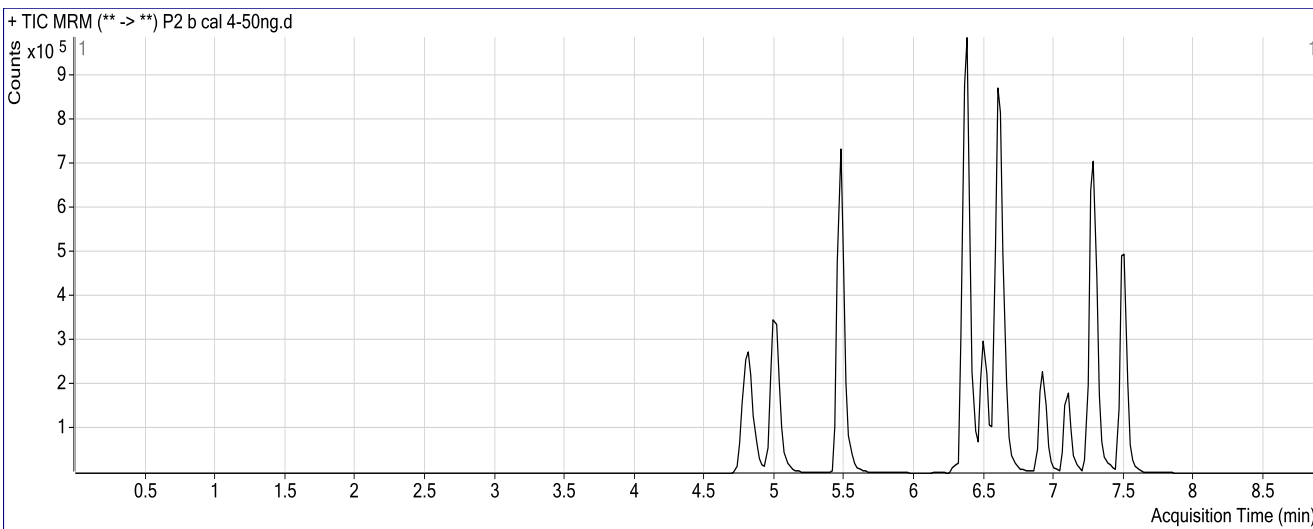
CJ

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 19:36 | <b>Data File</b>   | P2 b cal 4-50ng.d       |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 4-50ng.d       |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-D5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 372738   | 304007    | 1.2261     | 46.5251    |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 664119   | 161414    | 4.1144     | 46.4772    |



# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

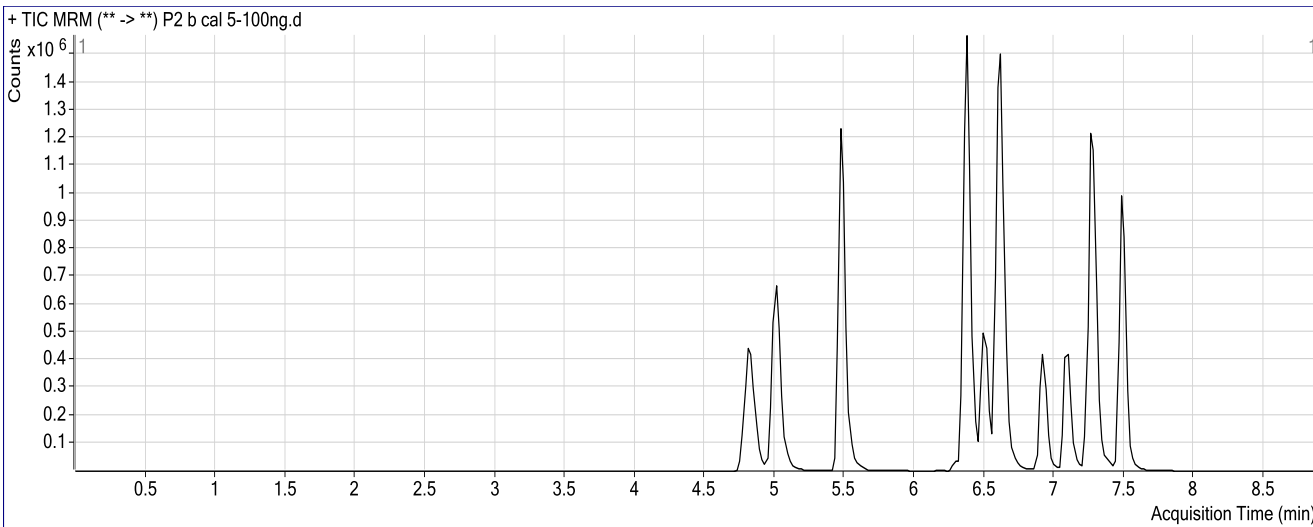
CJ

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 20:01 | <b>Data File</b>   | P2 b cal 5-100ng.d      |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 5-100ng.d      |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-E5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 800640   | 322681    | 2.4812     | 92.9661    |
| Flunitrazepam | Flunitrazepam-D7 | 7.477 | 1200780  | 140738    | 8.5321     | 95.3185    |

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

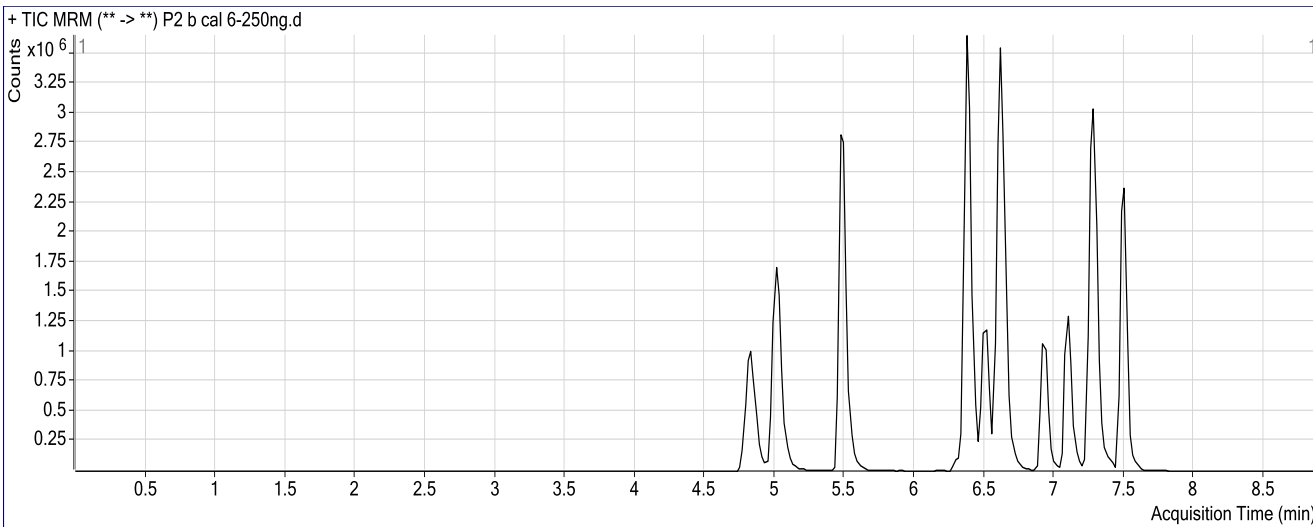
CS

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

### Analysis Info

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 20:27 | <b>Data File</b>   | P2 b cal 6-250ng.d      |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 6-250ng.d      |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-F5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

### Sample Chromatogram



### Results

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 2490623  | 354360    | 7.0285     | 261.2205   |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 2828557  | 126653    | 22.3331    | 247.9013   |

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

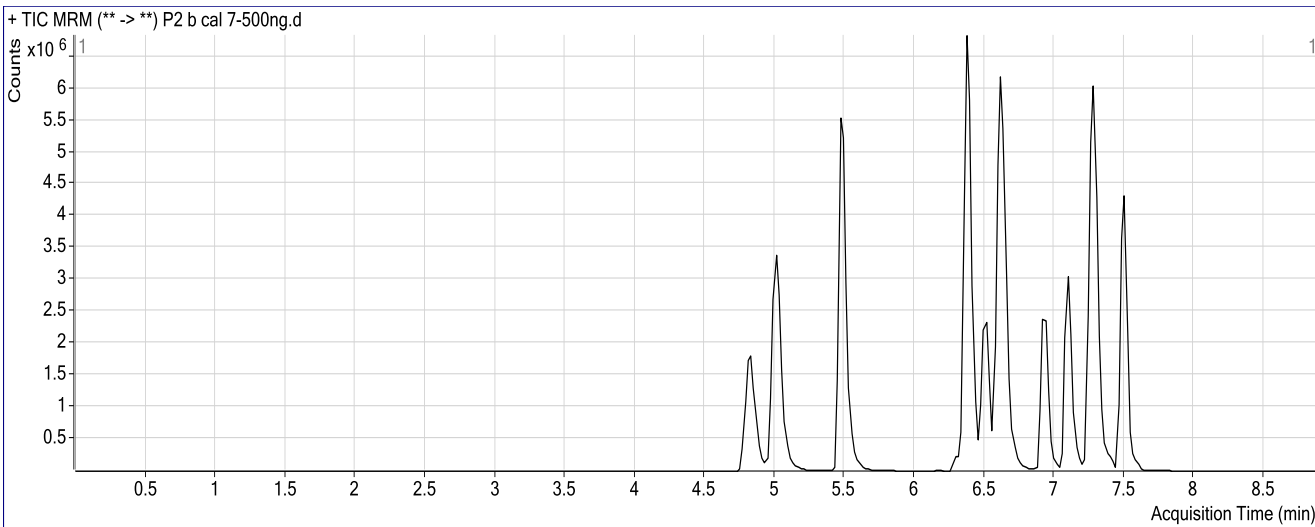
CJ

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 20:53 | <b>Data File</b>   | P2 b cal 7-500ng.d      |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 7-500ng.d      |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-G5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 5943604  | 416591    | 14.2672    | 529.0610   |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 4599576  | 90699     | 50.7126    | 561.6619   |

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

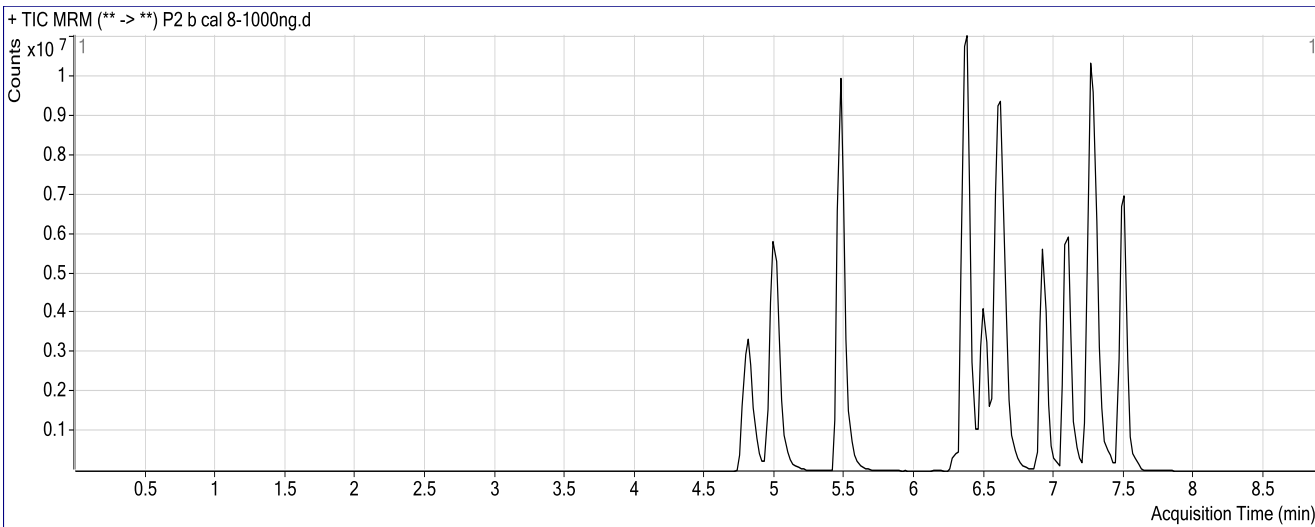
CJ

|                          |   |                      |           |
|--------------------------|---|----------------------|-----------|
| <b>Batch Data Path</b>   | C:\MassHunter\Data\2018\MD Quant\110218 MDQ P1 P2 CS\QuantResults\MDQ P2b 110218 CS.batch.bin |                      |           |
| <b>Analysis Time</b>     | 11/8/2018 2:09 PM   | <b>Analyst Name</b>  | ISPUser   |
| <b>Report Time</b>       | 11/9/2018 12:36 PM  | <b>Reporter Name</b> | ISPUser   |
| <b>Last Calib Update</b> | 11/8/2018 2:09 PM   | <b>Batch State</b>   | Processed |

**Analysis Info**

|                    |                  |                    |                         |
|--------------------|------------------|--------------------|-------------------------|
| <b>Acq Time</b>    | 2018-11-03 21:18 | <b>Data File</b>   | P2 b cal 8-1000ng.d     |
| <b>Sample Type</b> | Calibration      | <b>Sample Name</b> | P2 b cal 8-1000ng.d     |
| <b>Dilution</b>    | 1                | <b>Acq Method</b>  | MDQ Panel 2b 06-22-18.m |
| <b>Position</b>    | P1-H5            | <b>Sample Info</b> |                         |
| <b>Inj Vol</b>     | -1               | <b>Comment</b>     | AM 28-Panel 2           |

**Sample Chromatogram**



**Results**

| Compound      | ISTD Compound    | RT    | Response | ISTD Resp | Resp Ratio | Final Conc |
|---------------|------------------|-------|----------|-----------|------------|------------|
| Phencyclidine | Phencyclidine-D5 | 6.927 | 12685672 | 413116    | 30.7073    | 1137.3599  |
| Flunitrazepam | Flunitrazepam-D7 | 7.497 | 5740259  | 60284     | 95.2210    | 1053.7417  |