

D

**Worklist: 2638**

<u>LAB_CASE</u>	<u>ITEM</u>	<u>TASK_ID</u>	<u>DESCRIPTION</u>
M2018-3378	1	124341	AM 28 Blood Multi-Drug Quant Panel 2 by LC-QC
P2018-1953	1	124094	AM 28 Blood Multi-Drug Quant Panel 2 by LC-QC



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

B

Extraction Date: 08/15/18

Analyst: Sarah Pickle

Plate lot#: 0530563

Plate Expiration: 05/21/19

**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: 081518 P1 P2 SP Batch Name: *081518 MDQ P2a compounds removed*
- 2. Make necessary changes to integration limits
- 3. Integration linear and  $R^2$  values  $\geq 0.98$  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: Only zopiclone, carbamazepine, midazolam, and chlordiazepoxide were evaluated in this run.  
Curve range limited for carbamazepine 5-500.

B



# Idaho State Police Forensic Services

B

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

**Analyst:** Sarah Pickle  
**Extraction Date:** 08/15/18  
**Worklist Number:** 2638

<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	5/21/19			
Negative Blood	361331-1		05/25/18		
Methanol External Control Solution	022018	02/20/19	02/20/18		
Blood External Control Solution	081518	02/20/19	08/15/18		
0.5M Ammonium Hydroxide	032918	09/29/18	03/29/18		
Formic Acid (LCMS Grade)	095180B		11/15/17		
Water (LCMS Grade)	181370		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	081518	Make fresh	08/15/18		
5mM Ammonium Formate + 0.01% Formic Acid in Water (Mobile Phase A)	081518		08/15/18		
0.01% Formic Acid in Methanol (Mobile Phase B)	081518		08/15/18		
Needle Rinse--75% LCMS MeOH in LCMS Water	081518		08/15/18		

**Methanol External Control Solution (Lot: 022018)**  
100 ul each 1 mg/ml. stock solution in 9700 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		

**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:	02/20/19	

**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	181370

**5mM Ammonium Formate + 0.01% Formic Acid in Water (Lot: 081518)**

<i>Component</i>	<i>Source</i>	<i>Source Lot Number</i>
Ammonium Formate	Arcos Organics	A0287238
LCMS Water	Fisher	181370
Formic Acid	Fisher	095180B
Prepared:	08/15/18	
Prepared By:	Sarah Pickle	

**0.01% Formic Acid in LCMS Methanol (Mobile Phase B) (Lot: 081518)**

<i>Component</i>	<i>Source</i>	<i>Source Lot Number</i>
Formic Acid	Fisher	095180B
MeOH (LCMS Grade)	Fisher	172517
Prepared:	08/15/18	
Prepared By:	Sarah Pickle	

**Needle Rinse (75% LCMS MeOH in LCMS Water) (Lot: 081518)**

<i>Component</i>	<i>Source</i>	<i>Source Lot Number</i>
MeOH (LCMS Grade)	Fisher	177145
Water (LCMS Grade)	Fisher	177528
Prepared:	08/15/18	
Prepared By:	Sarah Pickle	

R

# ISP FORENSICS - Pocatello Instrument # 59740

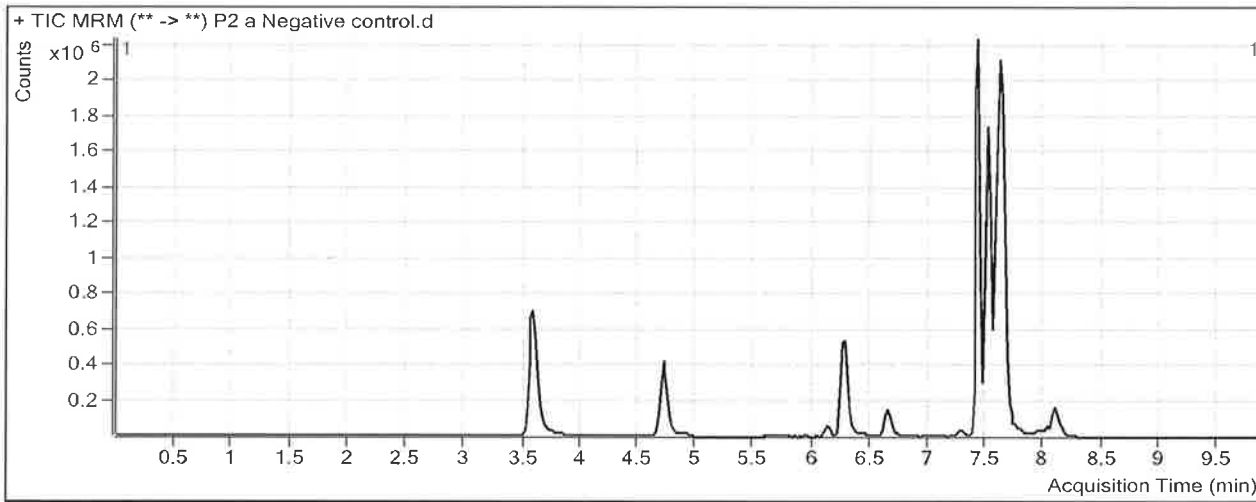
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

**Analysis Info**

**Acq Time** 2018-08-16 14:05 **Data File** P2 a Negative control.d  
**Sample Type** Sample **Sample Name** P2 a Negative control.d  
**Dilution** 1 **Acq Method** MDQ Panel 2a 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.131	1127	204420	0.0055	2.7755 < 5
Midazolam	Midazolam-D4	7.585	5320	2289121	0.0023	1.1992 < 5

B

# ISP FORENSICS - Pocatello Instrument # 59740

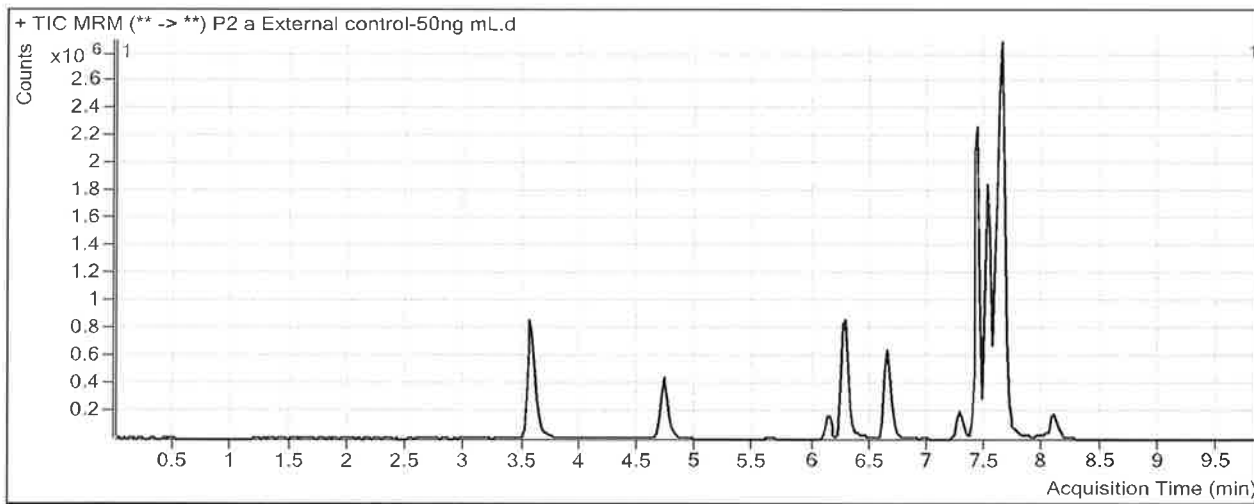
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 14:30 **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 06-22-18.m  
**Position** P2-F6 **Sample Info**  
**Inj Vol** -1 **Comment** AM 28; 361331-1 + WS022018 50 ng flunitrazepam, zopiclone, chlorodia:

### Sample Chromatogram



### Results

Compound	ISTD Compound	RT	Response	ISTD Resp	Resp Ratio	Final Conc
Zopiclone	Zopiclone-D4	6.131	339274	228761	1.4831	42.4431
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	1263622	4312530	0.2930	60.2573

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

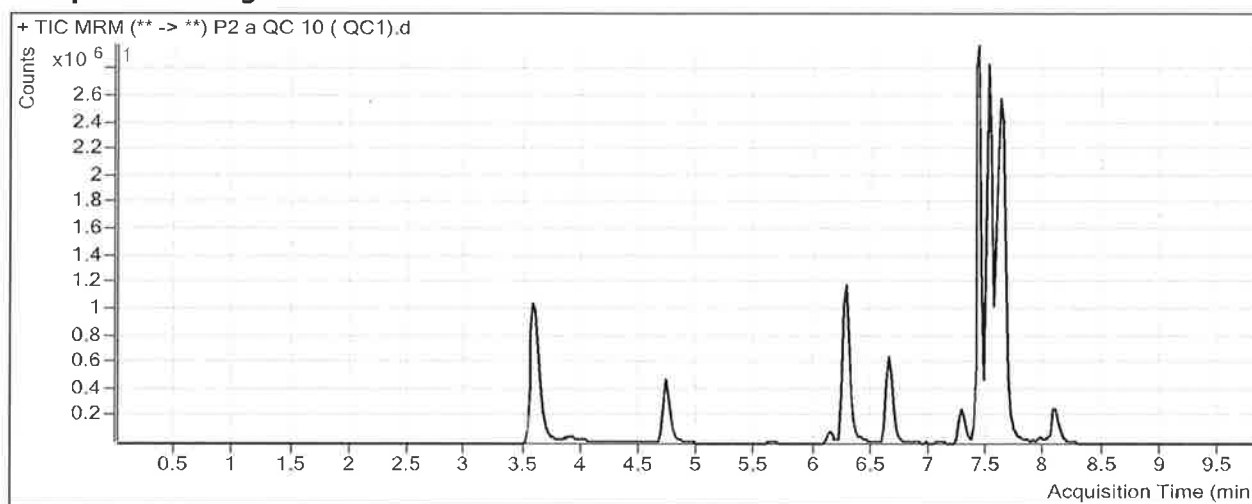
P

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 15:21 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.131	73623	235867	0.3121	11.0072
Carbamazepine	Carbamazepine-13C6	7.437	1819901	7315590	0.2488	9.9078
Midazolam	Midazolam-D4	7.585	153294	2709951	0.0566	9.1212
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	199649	4860486	0.0411	9.5636

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

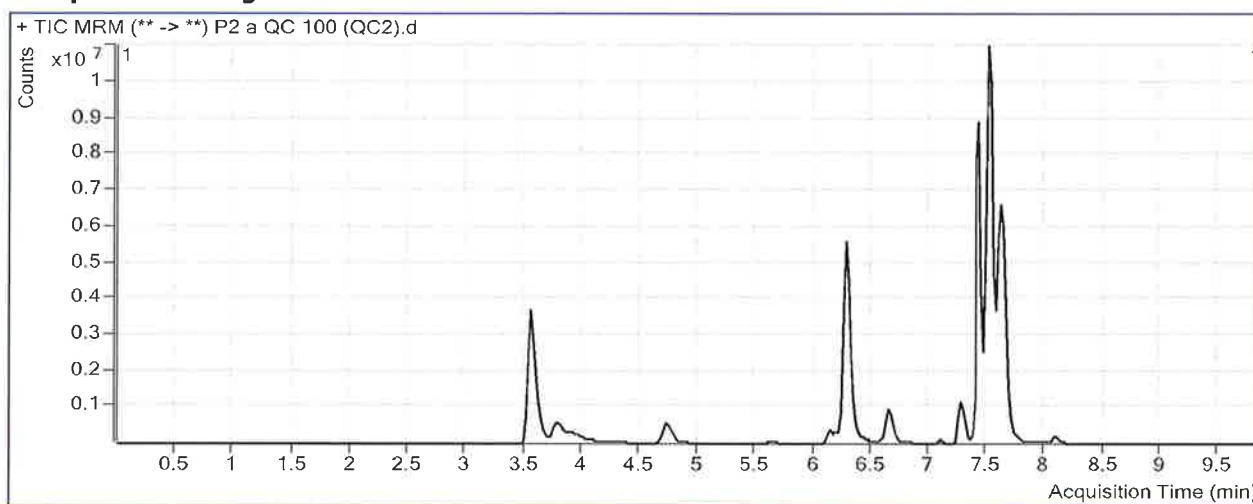
D

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:23 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 16:25 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.131	848182	254048	3.3387	92.2584
Carbamazepine	Carbamazepine-13C6	7.437	18370446	6907326	2.6596	110.1190
Midazolam	Midazolam-D4	7.585	2113257	2939868	0.7188	105.8430
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	1917750	4096622	0.4681	95.4940



# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

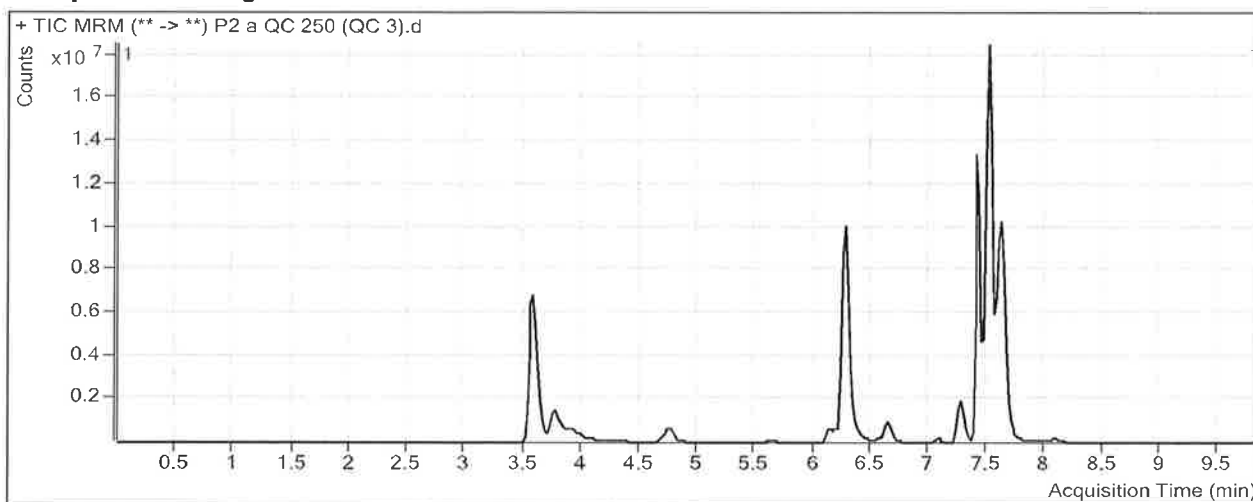
D

<b>Batch Data Path</b>	C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t		
<b>Analysis Time</b>	8/20/2018 3:20 PM	<b>Analyst Name</b>	ISPUser
<b>Report Time</b>	8/20/2018 3:23 PM	<b>Reporter Name</b>	ISPUser
<b>Last Calib Update</b>	8/20/2018 3:20 PM	<b>Batch State</b>	Processed

### Analysis Info

<b>Acq Time</b>	2018-08-16 17:28	<b>Data File</b>	P2 a QC 250 (QC 3).d
<b>Sample Type</b>	QC	<b>Sample Name</b>	P2 a QC 250 (QC 3).d
<b>Dilution</b>	1	<b>Acq Method</b>	MDQ Panel 2a 06-22-18.m
<b>Position</b>	P2-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
Zopiclone	Zopiclone-D4	6.111	1766244	200805	8.7958	238.7629
Carbamazepine	Carbamazepine-13C6	7.417	33513454	4759442	7.0415	292.2653
Midazolam	Midazolam-D4	7.565	4290171	2248003	1.9084	279.5828
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	3791199	2868278	1.3218	267.2605

SP

# ISP FORENSICS - Pocatello Instrument # 59740

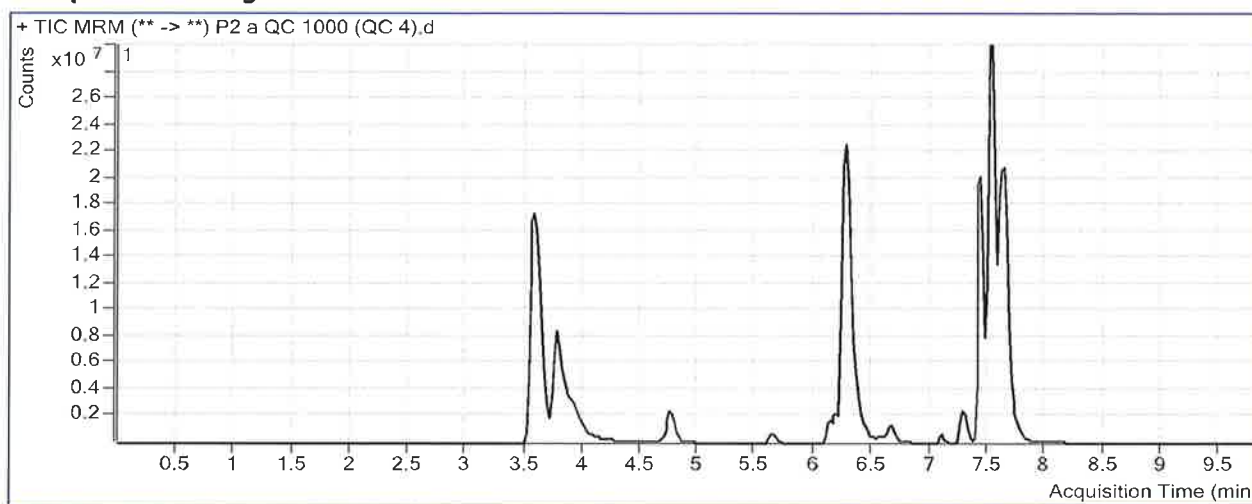
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:23 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 18:19 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.131	4511460	116983	38.5650	1037.9559
<del>Carbamazepine</del>	<del>Carbamazepine-13C6</del>	<del>7.437</del>	<del>65705814</del>	<del>4066415</del>	<del>16.1582</del>	<del>671.2264</del> *
Midazolam	Midazolam-D4	7.585	14361982	1945539	7.3820	1078.9860
Chlordiazepoxide	Chlordiazepoxide-D5	7.648	9763630	1974949	4.9437	996.0620

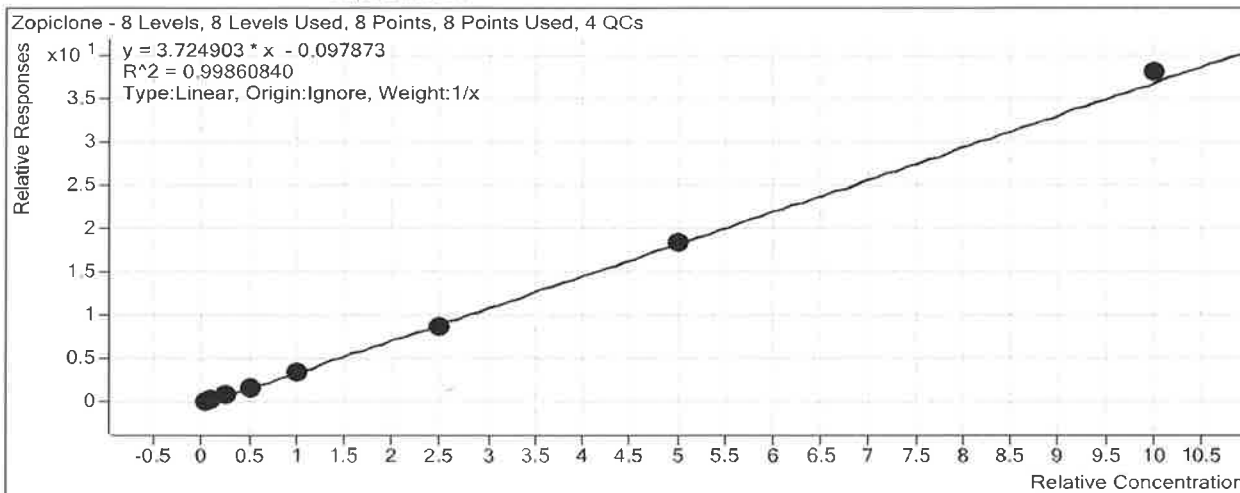
\*outside curve range. p

# ISP Forensics Calibration Curve Report



**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.batch.bin  
**Last Calib Update** 8/20/2018 3:20 PM **Analyst Name** ISP TOX

**Target Compound** Zopiclone  
**Internal Standard** Zopiclone-D4



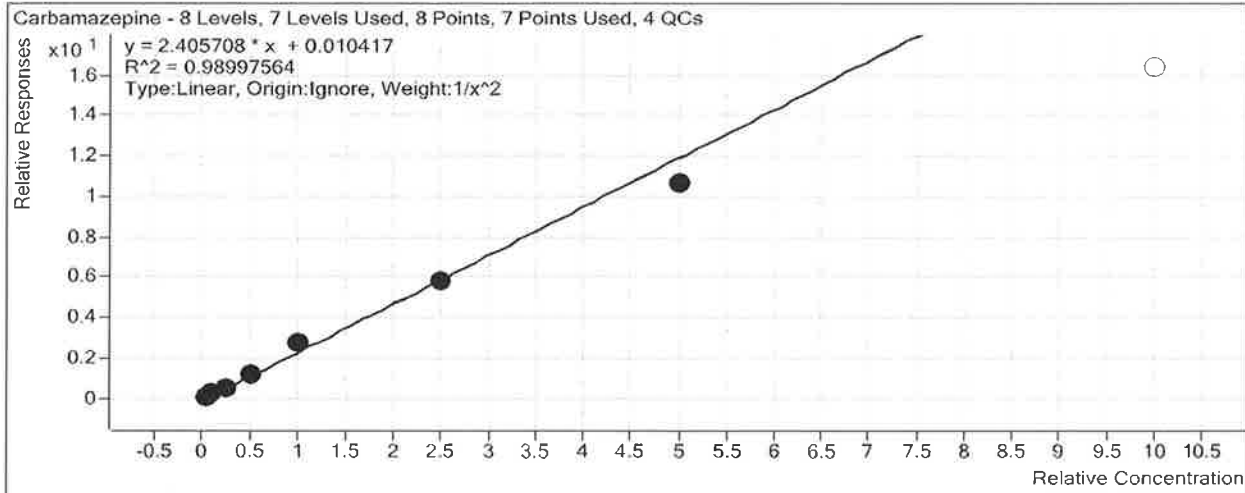
Sample	Level	Enabled	Exp Conc	Final Conc	Accuracy
P2 a cal 1-5ng.d	1	<input checked="" type="checkbox"/>	5	6.0	119.7
P2 a cal 2-10ng.d	2	<input checked="" type="checkbox"/>	10	10.7	107.3
P2 a QC 10 ( QC1).d	2	<input checked="" type="checkbox"/>	10	11.0	110.1
P2 a cal 3-25ng.d	3	<input checked="" type="checkbox"/>	25	22.3	89.1
P2 a cal 4-50ng.d	4	<input checked="" type="checkbox"/>	50	45.4	90.8
P2 a cal 5-100ng.d	5	<input checked="" type="checkbox"/>	100	96.5	96.5
P2 a QC 100 (QC2).d	5	<input checked="" type="checkbox"/>	100	92.3	92.3
P2 a cal 6-250ng.d	6	<input checked="" type="checkbox"/>	250	237.4	95.0
P2 a QC 250 (QC 3).d	6	<input checked="" type="checkbox"/>	250	238.8	95.5
P2 a cal 7-500ng.d	7	<input checked="" type="checkbox"/>	500	495.2	99.0
P2 a cal 8-1000ng.d	8	<input checked="" type="checkbox"/>	1000	1026.5	102.6
P2 a QC 1000 (QC 4).d	8	<input checked="" type="checkbox"/>	1000	1038.0	103.8

# ISP Forensics Calibration Curve Report

D

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.batch.bin  
**Last Calib Update** 8/20/2018 3:20 PM **Analyst Name** ISP TOX

**Target Compound** Carbamazepine  
**Internal Standard** Carbamazepine-13C6



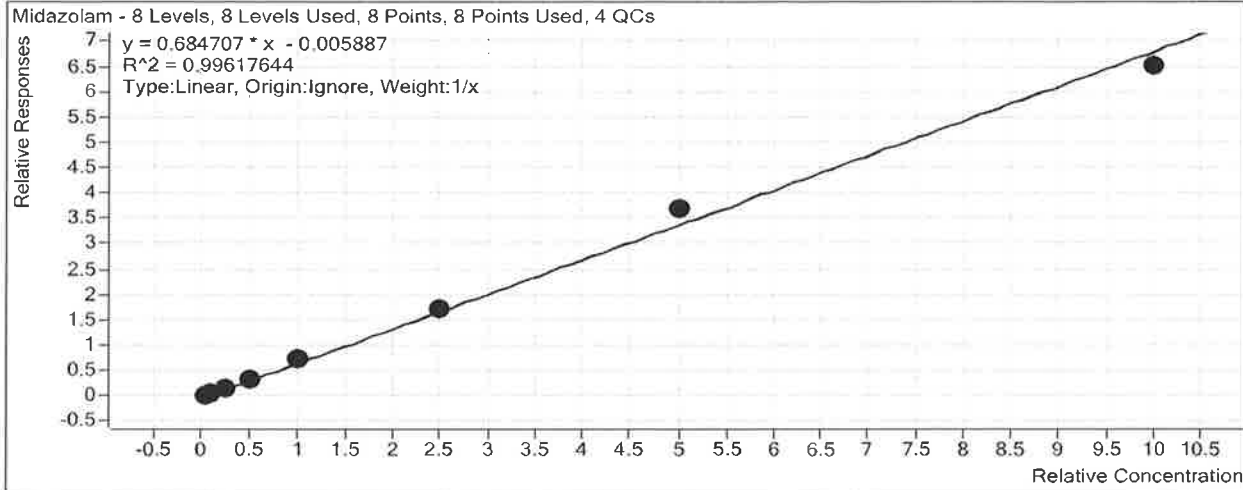
Sample	Level	Enabled	Exp Conc	Final Conc	Accuracy
P2 a cal 1-5ng.d	1	<input checked="" type="checkbox"/>	5	4.9	97.0
P2 a cal 2-10ng.d	2	<input checked="" type="checkbox"/>	10	10.6	106.4
P2 a QC 10 ( QC1).d	2	<input checked="" type="checkbox"/>	10	9.9	99.1
P2 a cal 3-25ng.d	3	<input checked="" type="checkbox"/>	25	23.9	95.6
P2 a cal 4-50ng.d	4	<input checked="" type="checkbox"/>	50	50.5	101.1
P2 a cal 5-100ng.d	5	<input checked="" type="checkbox"/>	100	114.8	114.8
P2 a QC 100 (QC2).d	5	<input checked="" type="checkbox"/>	100	110.1	110.1
P2 a cal 6-250ng.d	6	<input checked="" type="checkbox"/>	250	241.9	96.7
P2 a QC 250 (QC 3).d	6	<input checked="" type="checkbox"/>	250	292.3	116.9
P2 a cal 7-500ng.d	7	<input checked="" type="checkbox"/>	500	441.9	88.4
P2 a cal 8-1000ng.d	8	<input type="checkbox"/>	1000	680.1	68.0
P2 a QC 1000 (QC 4).d	8	<input type="checkbox"/>	1000	671.2	67.1

# ISP Forensics Calibration Curve Report

D

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.batch.bin  
**Last Calib Update** 8/20/2018 3:20 PM **Analyst Name** ISP TOX

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



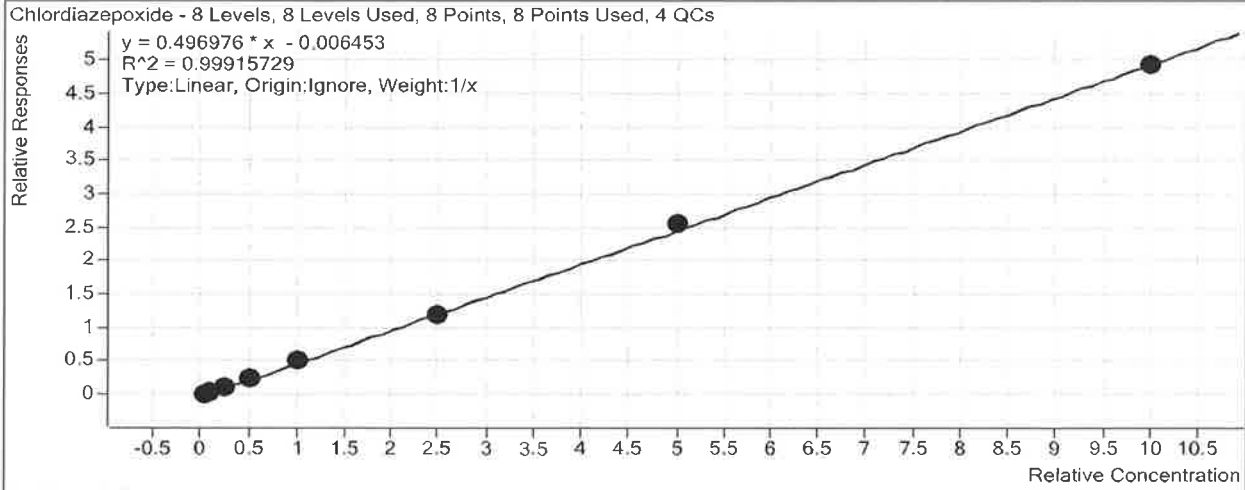
Sample	Level	Enabled	Exp Conc	Final Conc	Accuracy
P2 a cal 1-5ng.d	1	<input checked="" type="checkbox"/>	5	5.1	101.6
P2 a cal 2-10ng.d	2	<input checked="" type="checkbox"/>	10	9.9	99.2
P2 a QC 10 ( QC1).d	2	<input checked="" type="checkbox"/>	10	9.1	91.2
P2 a cal 3-25ng.d	3	<input checked="" type="checkbox"/>	25	21.5	86.2
P2 a cal 4-50ng.d	4	<input checked="" type="checkbox"/>	50	50.7	101.4
P2 a cal 5-100ng.d	5	<input checked="" type="checkbox"/>	100	106.9	106.9
P2 a QC 100 (QC2).d	5	<input checked="" type="checkbox"/>	100	105.8	105.8
P2 a cal 6-250ng.d	6	<input checked="" type="checkbox"/>	250	253.6	101.4
P2 a QC 250 (QC 3).d	6	<input checked="" type="checkbox"/>	250	279.6	111.8
P2 a cal 7-500ng.d	7	<input checked="" type="checkbox"/>	500	541.2	108.2
P2 a cal 8-1000ng.d	8	<input checked="" type="checkbox"/>	1000	951.1	95.1
P2 a QC 1000 (QC 4).d	8	<input checked="" type="checkbox"/>	1000	1079.0	107.9

# ISP Forensics Calibration Curve Report

P

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.batch.bin  
**Last Calib Update** 8/20/2018 3:20 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	5.2	104.3
P2 a cal 2-10ng.d	2	<input checked="" type="checkbox"/>	10	10.0	100.1
P2 a QC 10 ( QC1).d	2	<input checked="" type="checkbox"/>	10	9.6	95.6
P2 a cal 3-25ng.d	3	<input checked="" type="checkbox"/>	25	22.9	91.8
P2 a cal 4-50ng.d	4	<input checked="" type="checkbox"/>	50	50.1	100.1
P2 a cal 5-100ng.d	5	<input checked="" type="checkbox"/>	100	105.8	105.8
P2 a QC 100 (QC2).d	5	<input checked="" type="checkbox"/>	100	95.5	95.5
P2 a cal 6-250ng.d	6	<input checked="" type="checkbox"/>	250	239.2	95.7
P2 a QC 250 (QC 3).d	6	<input checked="" type="checkbox"/>	250	267.3	106.9
P2 a cal 7-500ng.d	7	<input checked="" type="checkbox"/>	500	515.0	103.0
P2 a cal 8-1000ng.d	8	<input checked="" type="checkbox"/>	1000	991.7	99.2
P2 a QC 1000 (QC 4).d	8	<input checked="" type="checkbox"/>	1000	996.1	99.6

# ISP FORENSICS - Pocatello Instrument # 59740

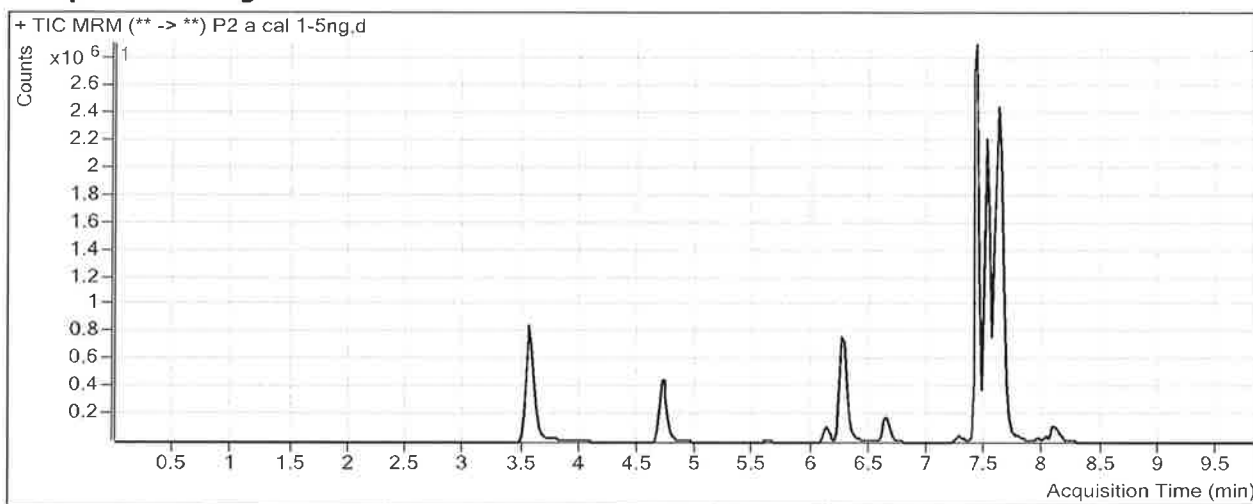
## Multi-Drug Confirmatory Analysis Report

<b>Batch Data Path</b>	C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t		
<b>Analysis Time</b>	8/20/2018 3:20 PM	<b>Analyst Name</b>	ISPUser
<b>Report Time</b>	8/20/2018 3:21 PM	<b>Reporter Name</b>	ISPUser
<b>Last Calib Update</b>	8/20/2018 3:20 PM	<b>Batch State</b>	Processed

### Analysis Info

<b>Acq Time</b>	2018-08-16 10:15	<b>Data File</b>	P2 a cal 1-5ng.d
<b>Sample Type</b>	Calibration	<b>Sample Name</b>	P2 a cal 1-5ng.d
<b>Dilution</b>	1	<b>Acq Method</b>	MDQ Panel 2a 06-22-18.m
<b>Position</b>	P2-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
Zopiclone	Zopiclone-D4	6.111	49786	398350	0.1250	5.9828
Carbamazepine	Carbamazepine-13C6	7.417	1040375	8185551	0.1271	4.8502
Midazolam	Midazolam-D4	7.585	73491	2543605	0.0289	5.0794
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	94100	4832543	0.0195	5.2165

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

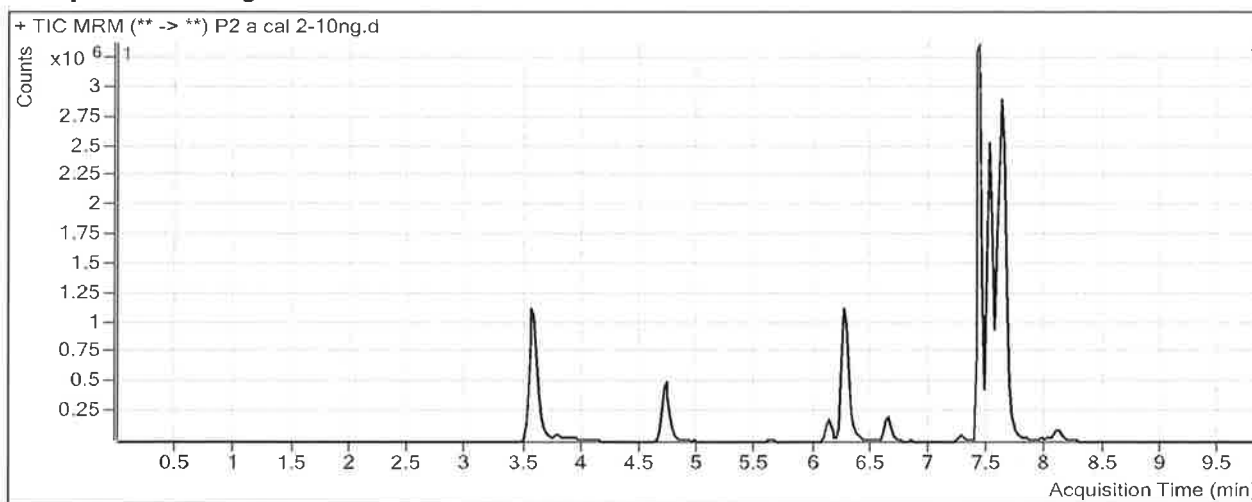
P

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:21 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 10:40 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	154292	511388	0.3017	10.7274
Carbamazepine	Carbamazepine-13C6	7.417	2317430	8698175	0.2664	10.6418
Midazolam	Midazolam-D4	7.585	174368	2810811	0.0620	9.9198
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	227574	5256218	0.0433	10.0103



# ISP FORENSICS - Pocatello Instrument # 59740

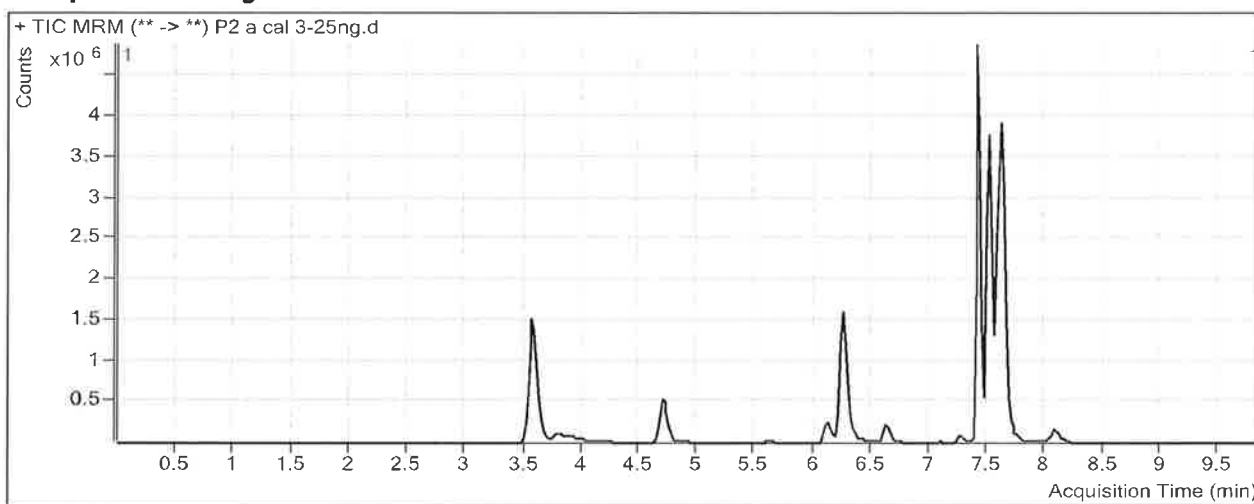
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:21 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 11:06 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	372960	509686	0.7317	22.2722
Carbamazepine	Carbamazepine-13C6	7.417	5647751	9646375	0.5855	23.9041
Midazolam	Midazolam-D4	7.585	438624	3097541	0.1416	21.5407
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	610307	5674097	0.1076	22.9413

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

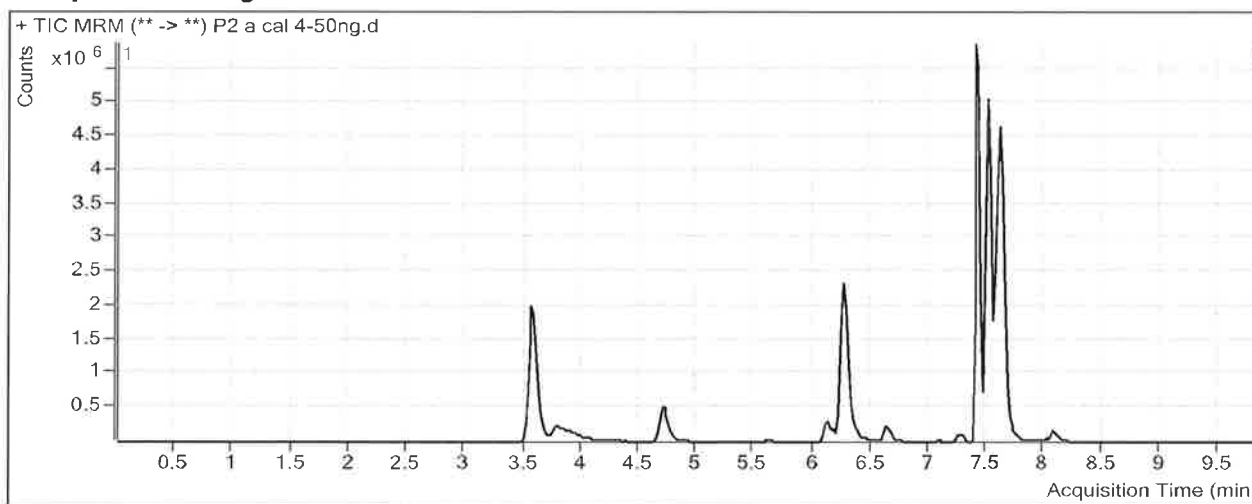
P

<b>Batch Data Path</b>	C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t		
<b>Analysis Time</b>	8/20/2018 3:20 PM	<b>Analyst Name</b>	ISPUser
<b>Report Time</b>	8/20/2018 3:22 PM	<b>Reporter Name</b>	ISPUser
<b>Last Calib Update</b>	8/20/2018 3:20 PM	<b>Batch State</b>	Processed

### Analysis Info

<b>Acq Time</b>	2018-08-16 11:31	<b>Data File</b>	P2 a cal 4-50ng.d
<b>Sample Type</b>	Calibration	<b>Sample Name</b>	P2 a cal 4-50ng.d
<b>Dilution</b>	1	<b>Acq Method</b>	MDQ Panel 2a 06-22-18.m
<b>Position</b>	P2-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
Zopiclone	Zopiclone-D4	6.111	594664	373081	1.5939	45.4187
Carbamazepine	Carbamazepine-13C6	7.417	10132143	8263937	1.2261	50.5319
Midazolam	Midazolam-D4	7.585	922436	2703723	0.3412	50.6873
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	1161730	4792468	0.2424	50.0749

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

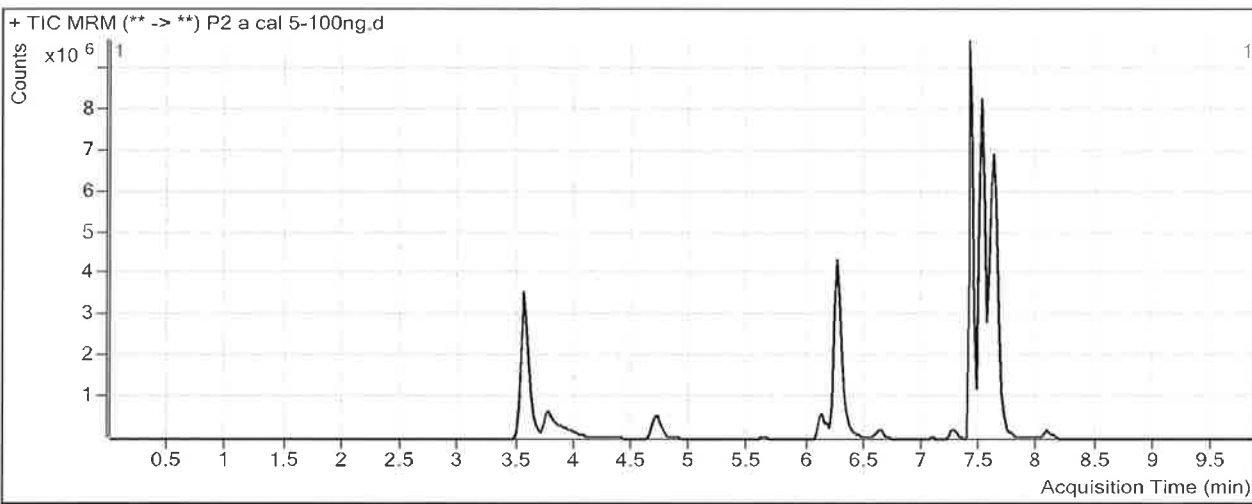
P

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 11:57 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	1395427	399139	3.4961	96.4849
Carbamazepine	Carbamazepine-13C6	7.417	22173941	8000899	2.7714	114.7693
Midazolam	Midazolam-D4	7.565	1983261	2731447	0.7261	106.9028
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	2326962	4481444	0.5192	105.7791

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

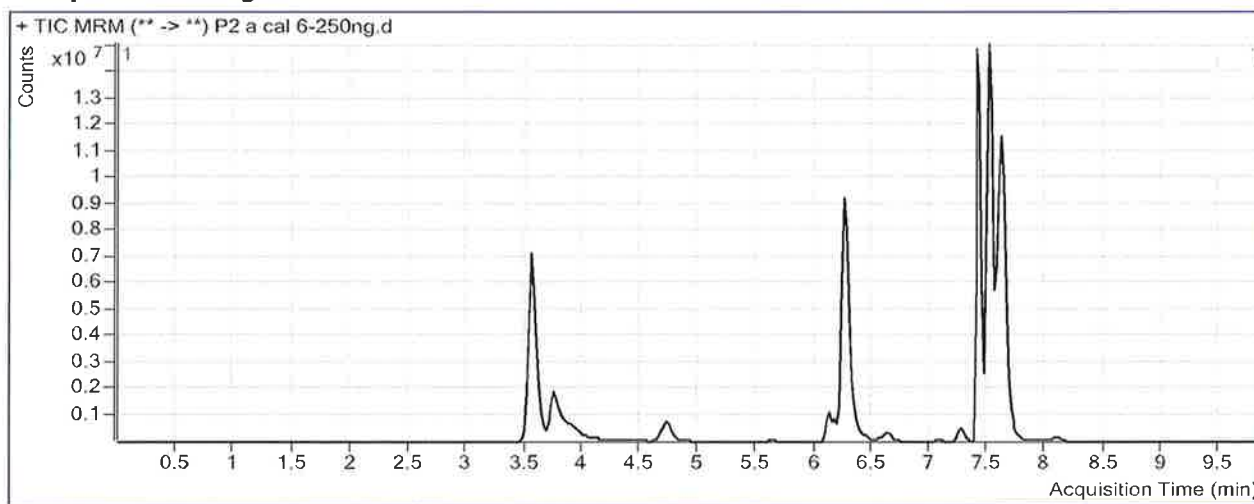
D

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 12:22 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	2751111	314552	8.7461	237.4287
Carbamazepine	Carbamazepine-13C6	7.417	41936571	7194724	5.8288	241.8572
Midazolam	Midazolam-D4	7.565	4310878	2491409	1.7303	253.5660
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	4621806	3908319	1.1826	239.2487

B

# ISP FORENSICS - Pocatello Instrument # 59740

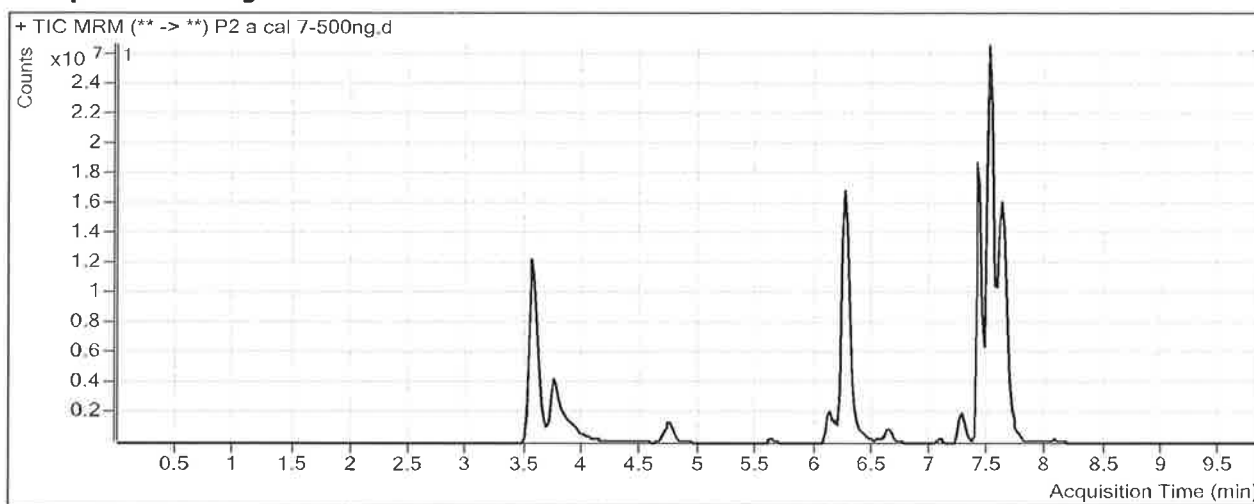
## Multi-Drug Confirmatory Analysis Report

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.L  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

### Analysis Info

**Acq Time** 2018-08-16 12:48 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	5501799	299862	18.3478	495.1979
Carbamazepine	Carbamazepine-13C6	7.417	58006069	5450714	10.6419	441.9282
Midazolam	Midazolam-D4	7.565	7964859	2152940	3.6995	541.1677
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	7383798	2892286	2.5529	514.9907

# ISP FORENSICS - Pocatello Instrument # 59740

## Multi-Drug Confirmatory Analysis Report

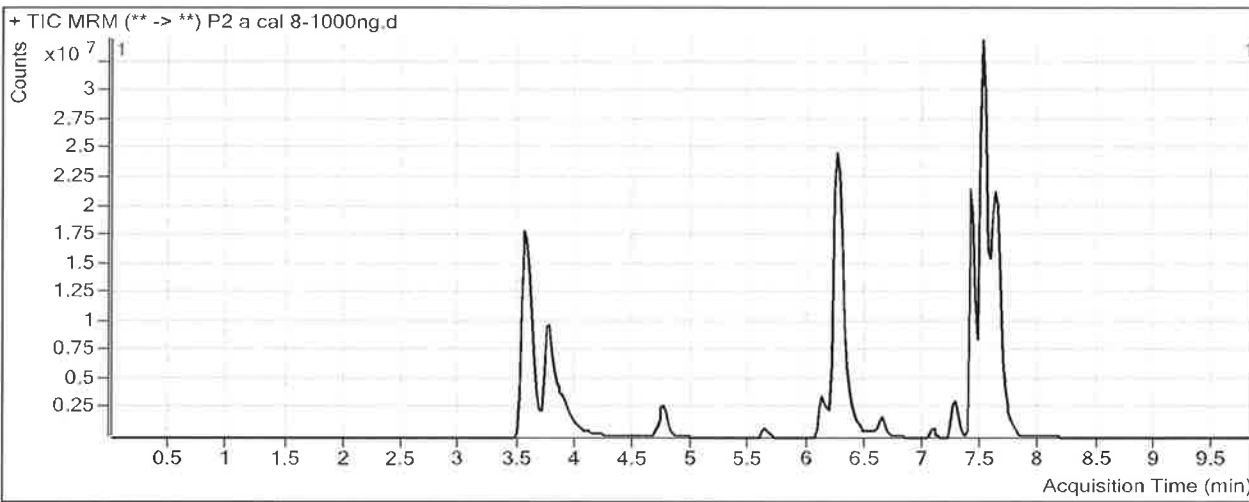
P

**Batch Data Path** C:\MassHunter\Data\2018\MD Quant\081518 P1 P2 SP\QuantResults\081518 MDQ P2a compounds removed.t  
**Analysis Time** 8/20/2018 3:20 PM **Analyst Name** ISPUser  
**Report Time** 8/20/2018 3:22 PM **Reporter Name** ISPUser  
**Last Calib Update** 8/20/2018 3:20 PM **Batch State** Processed

**Analysis Info**

**Acq Time** 2018-08-16 13:14 **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 06-22-18.m  
**Position** P2-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
Zopiclone	Zopiclone-D4	6.111	9276093	243226	38.1378	1026.4875
Carbamazepine	Carbamazepine-13C6	7.417	73470135	4487358	16.3727	680.1438
Midazolam	Midazolam-D4	7.565	11651140	1790662	6.5066	951.1363
Chlordiazepoxide	Chlordiazepoxide-D5	7.627	11190447	2273441	4.9223	991.7384