Quantitation Report (Qedit)

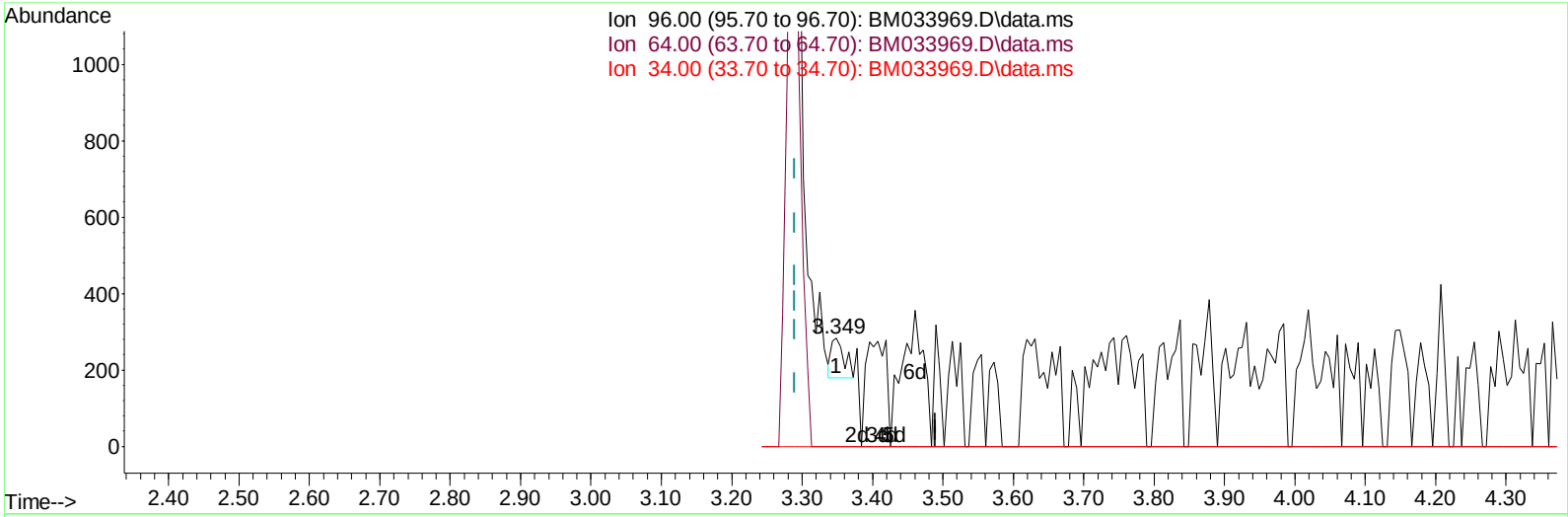
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033969.D
Acq On : 02 Feb 2022 11:45
Operator : CG/JU
Sample : N1273-21DL 5X
Misc :
ALS Vial : 40 Sample Multiplier: 1

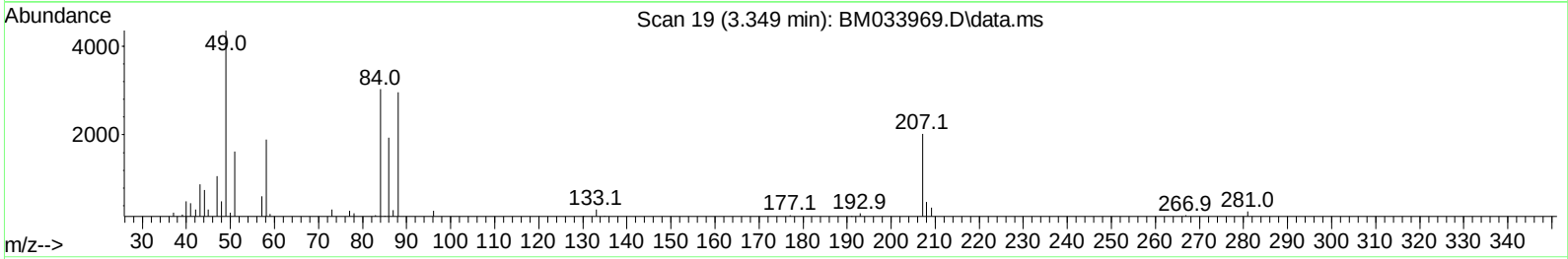
Quant Time: Feb 03 00:11:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

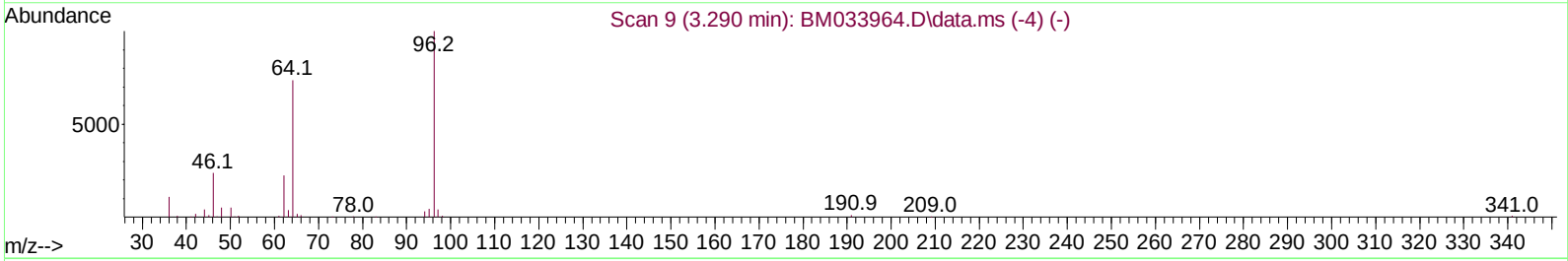
Instrument :
BNA_M
ClientSampleId :
C0HL9DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :Yogesh Patel 02/04/2022







TIC: BM033969.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.349min (+ 0.059) 0.05 ng/uL

response 131

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	0.00#
34.00	0.00	0.00
0.00	0.00	0.00

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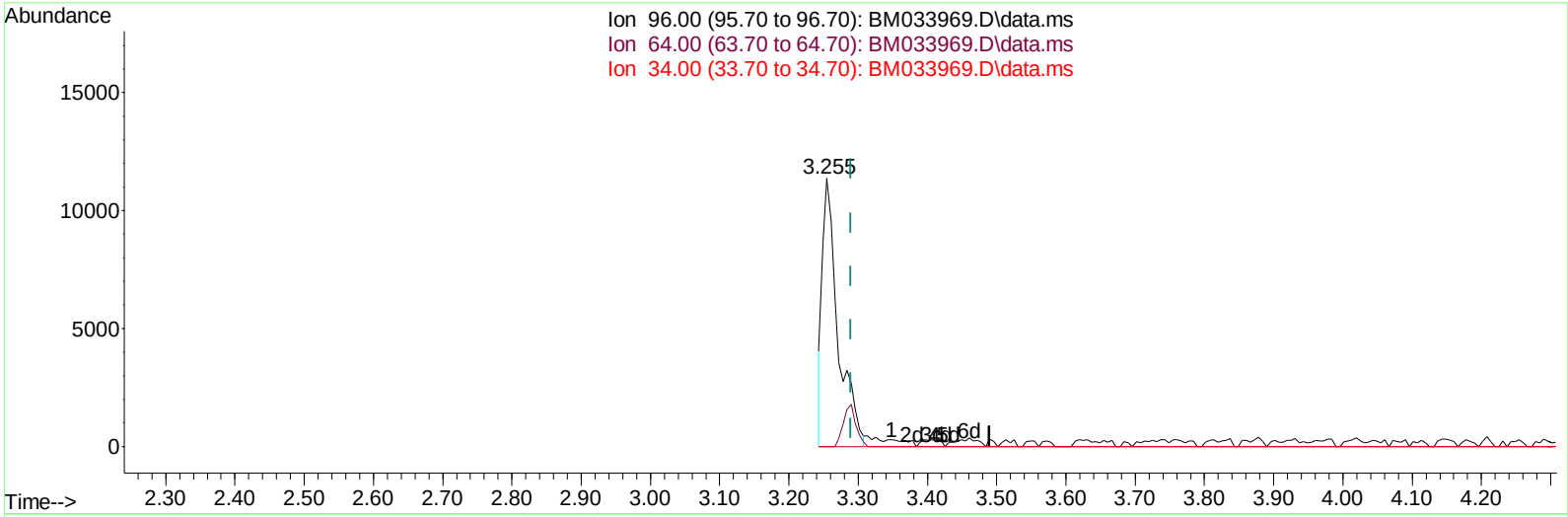

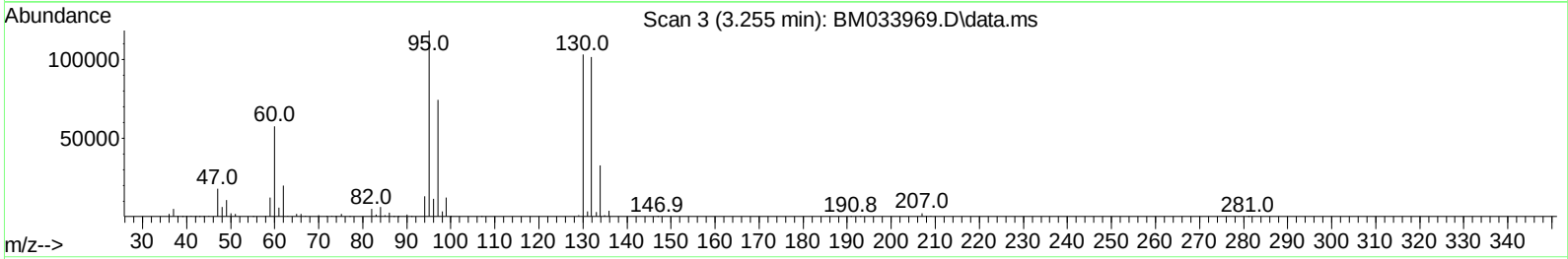

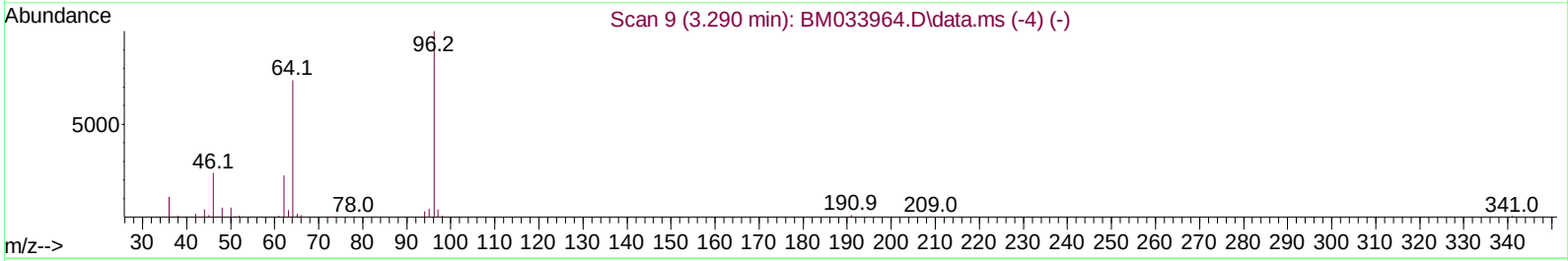

TIC: BM033969.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.255min (-0.035) 6.33 ng/uL m

response 17921

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	0.00#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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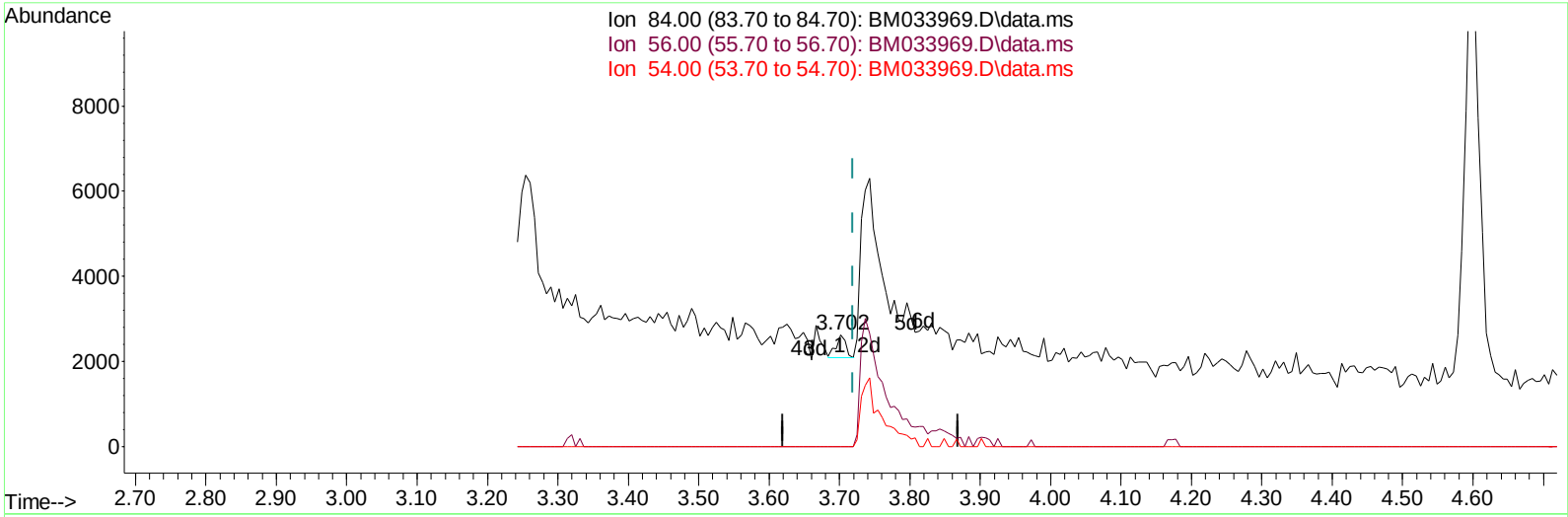
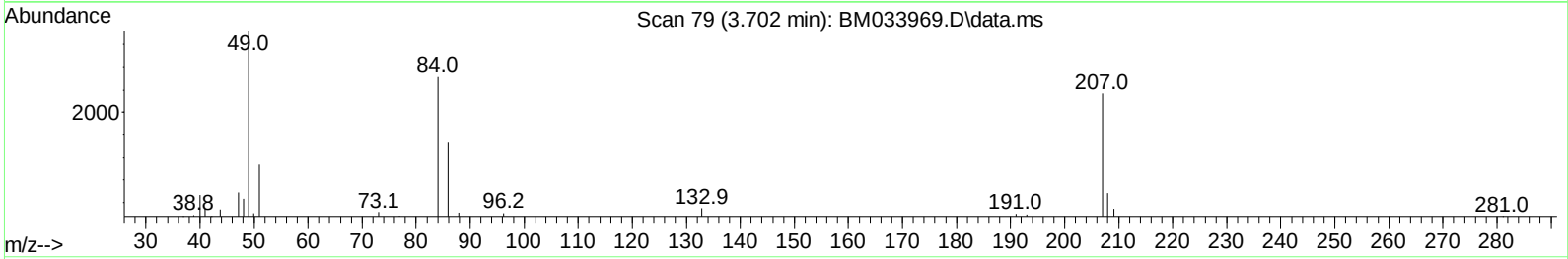
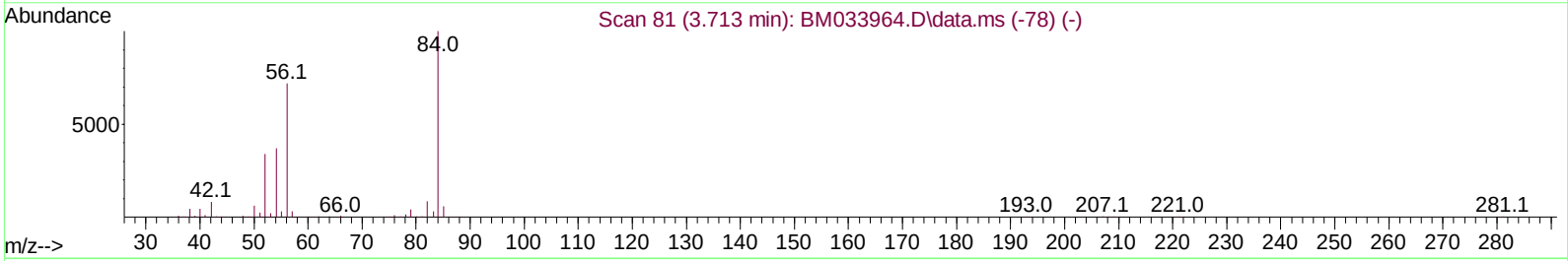
Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :Yogesh Patel 02/04/2022







TIC: BM033969.D\data.ms

(4) Pyridine-d5 (S)

3.702min (-0.018) 0.06 ng/ul

response 495

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	0.00#
54.00	42.60	0.00#
0.00	0.00	0.00

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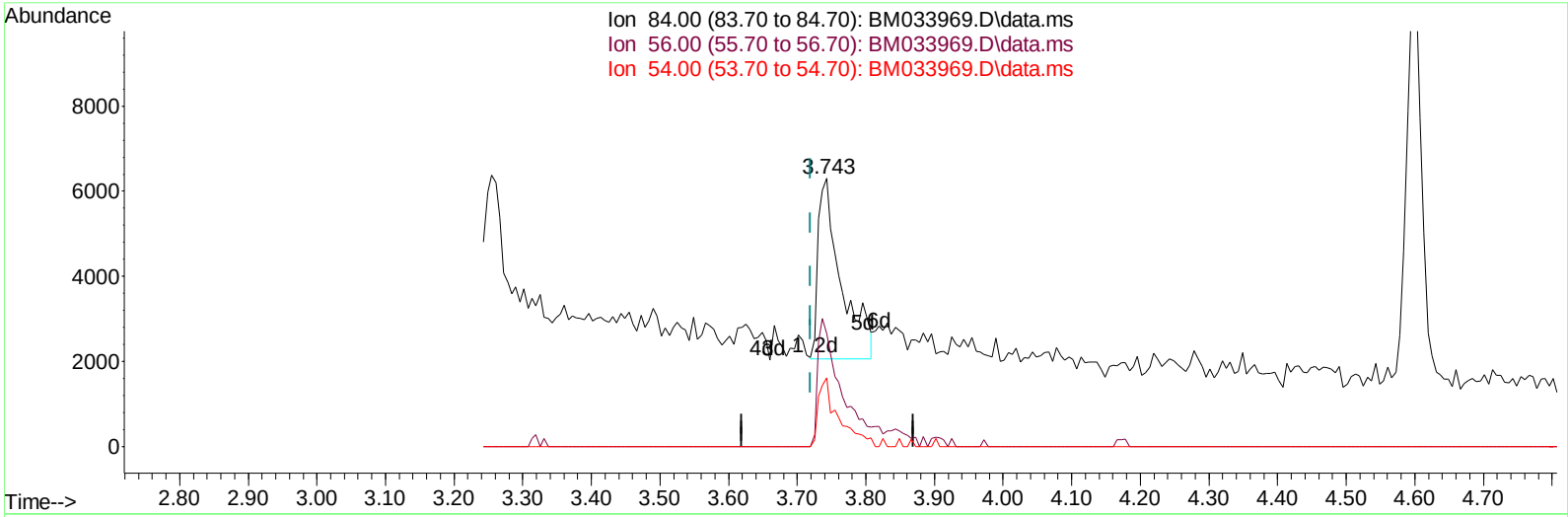
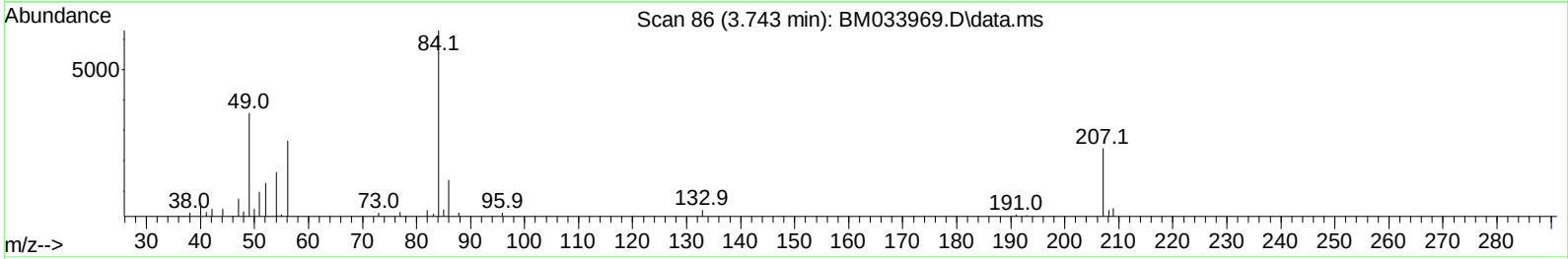
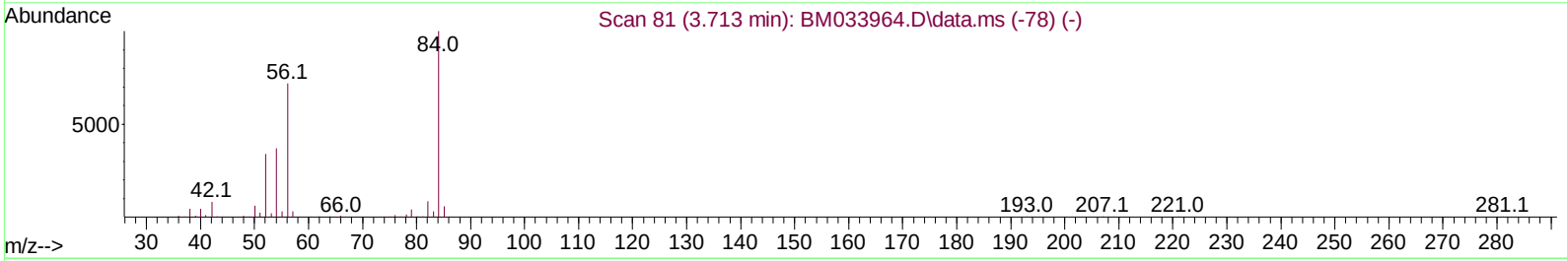
ClientSampleId :
C0HL9DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :Yogesh Patel 02/04/2022







TIC: BM033969.D\data.ms

(4) Pyridine-d5 (S)

3.743min (+ 0.024) 1.26 ng/ul m

response 9960

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	42.21#
54.00	42.60	25.59#
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/03/2022
Supervised By :Yogesh Patel 02/04/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	111742	20.000	ng/ul	0.00
20) Naphthalene-d8	10.736	136	488093	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	351507	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	753806	20.000	ng/ul	0.00
79) Chrysene-d12	21.459	240	594463	20.000	ng/ul	0.00
88) Perylene-d12	23.853	264	527498	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.255	96	17921m	6.325	ng/uL	-0.04
4) Pyridine-d5	3.743	84	9960m	1.259	ng/ul	0.02
7) Phenol-d5	7.084	99	10743	1.148	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	44719	7.836	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	36120	4.816	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	20406	2.751	ng/ul	0.00
21) Nitrobenzene-d5	9.084	128	26868	6.860	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143	23178	5.480	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165	38532	5.035	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	58795	5.563	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166	217272	8.272	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	230413	7.161	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.554	176	183032	8.345	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	19895	4.204	ng/ul	0.00
73) Anthracene-d10	17.395	188	293697	8.136	ng/ul	0.00
81) Pyrene-d10	19.683	212	320618	9.135	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.694	264	228822	8.259	ng/ul	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.319	88	74977	24.250	ng/uL#	90

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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