Quantitation Report (Qedit)

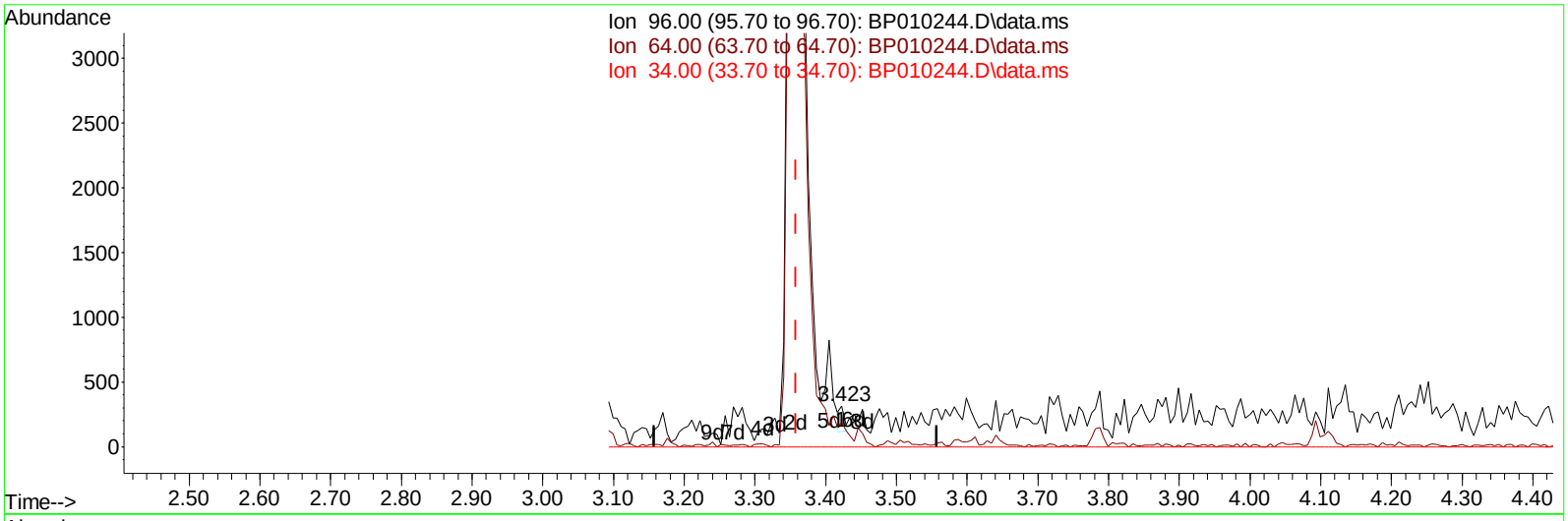
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050522\
Data File : BP010244.D
Acq On : 05 May 2022 13:29
Operator : CG/JU
Sample : N2668-14
Misc :
ALS Vial : 7 Sample Multiplier: 1

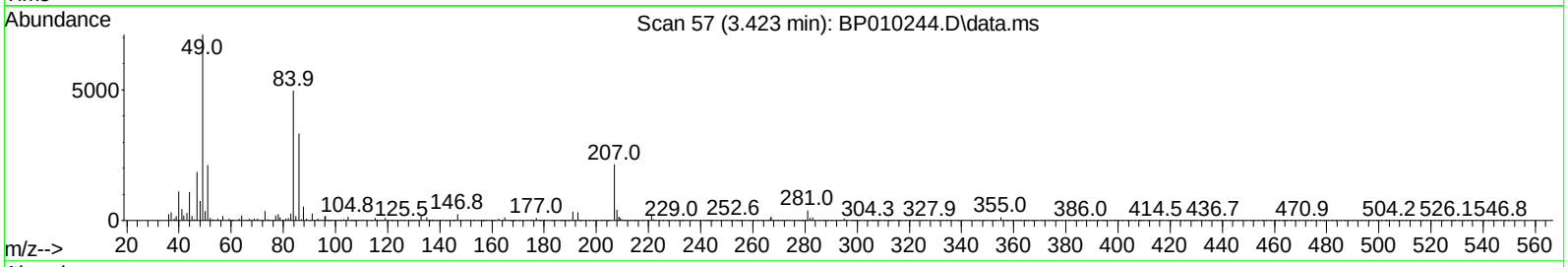
Quant Time: May 06 01:06:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 06 01:05:07 2022
Response via : Initial Calibration

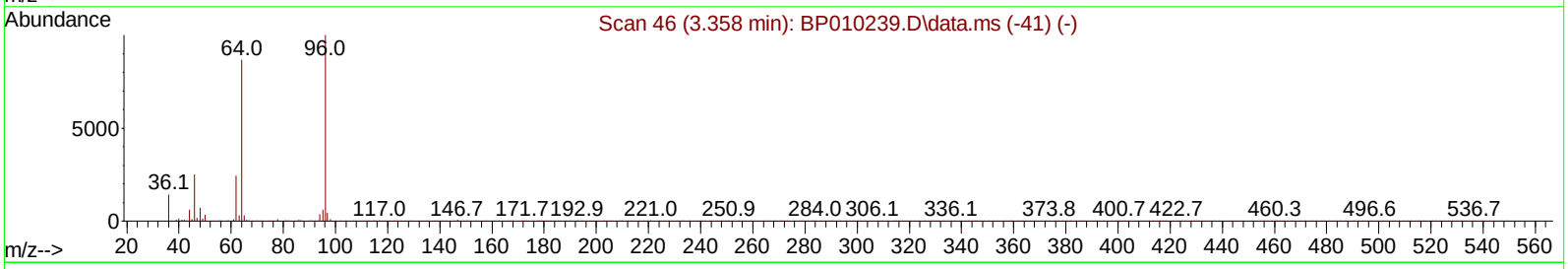
Instrument :
BNA_P
ClientSampleId :
EXL71

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/10/2022
Supervised By :mohammad ahmed 05/10/2022







TIC: BP010244.D\data.ms

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

3.423min (+ 0.065) 0.03 ng/uL

response 89

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.18
34.00	0.00	0.00
0.00	0.00	0.00

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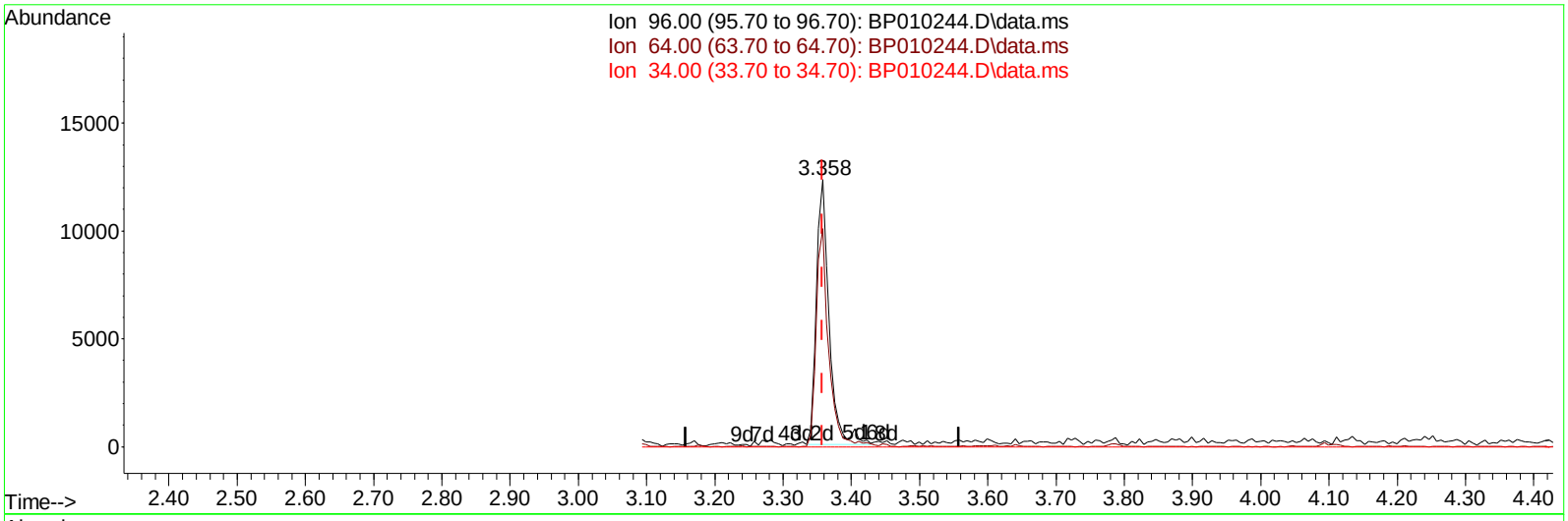

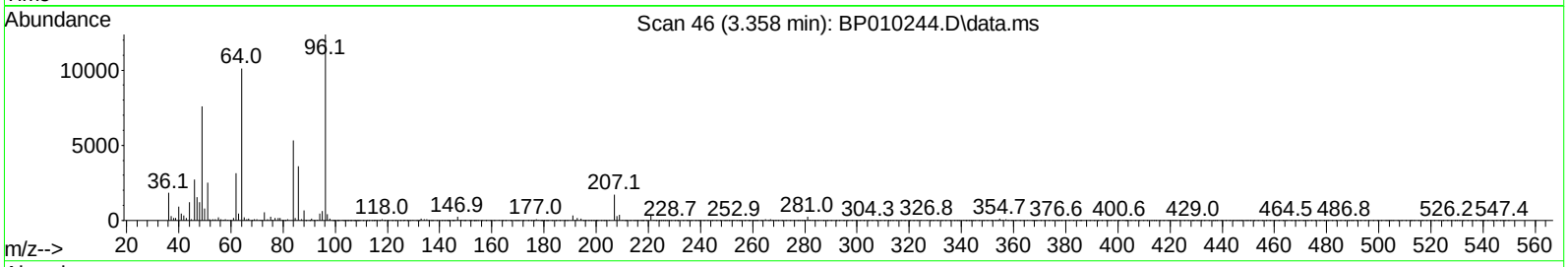

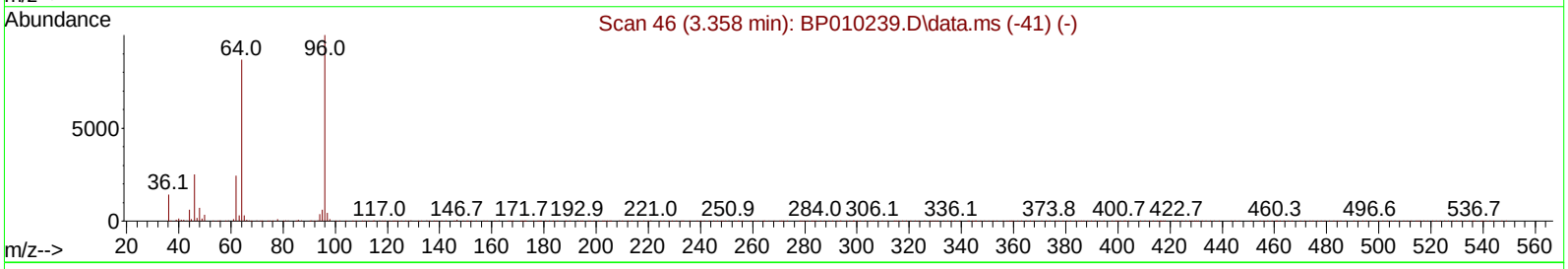

TIC: BP010244.D\data.ms

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

3.358min (+ 0.000) 4.51 ng/uL m

response 15962

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.71#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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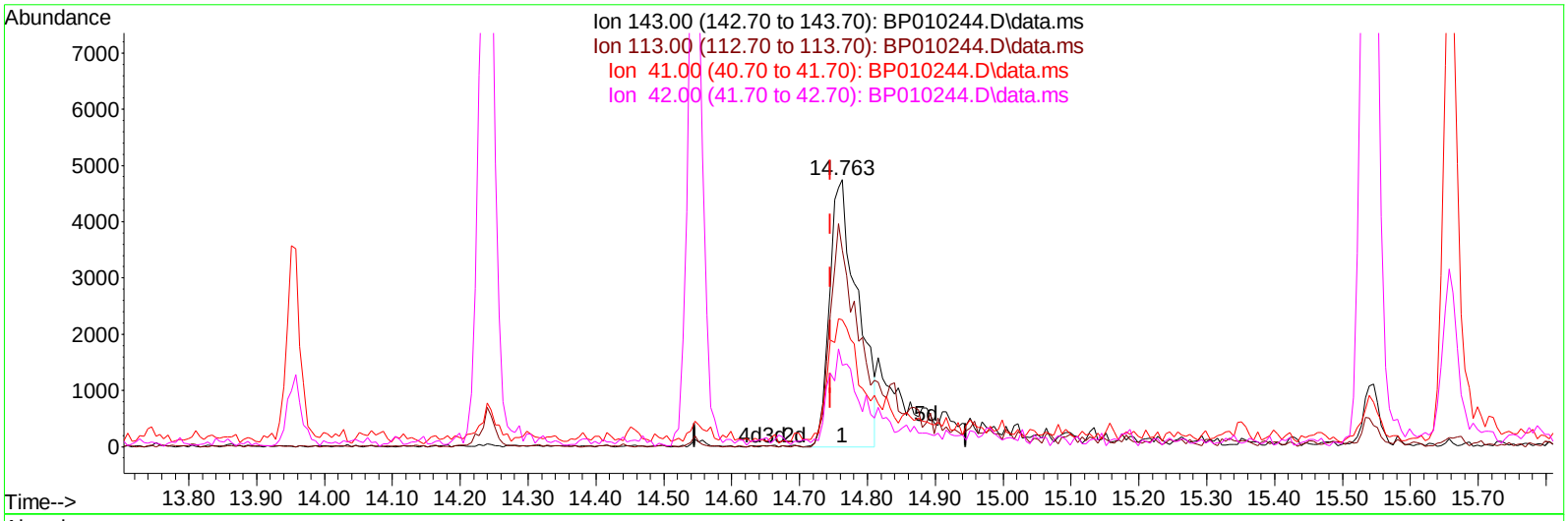

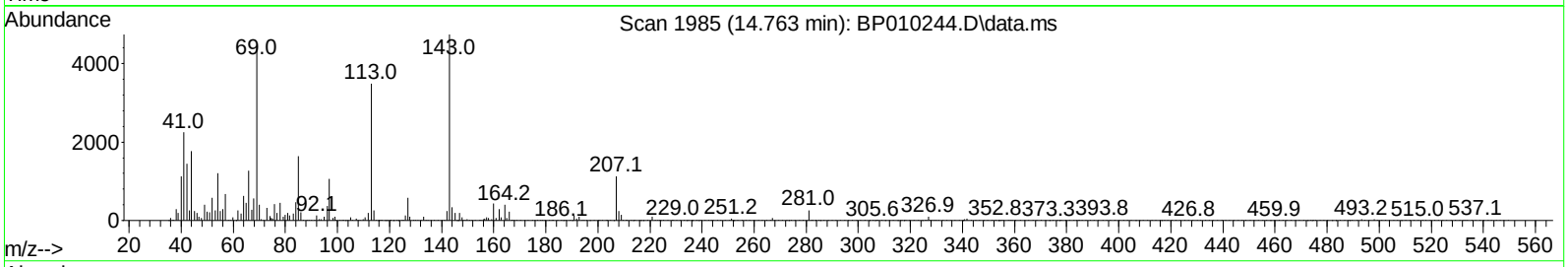

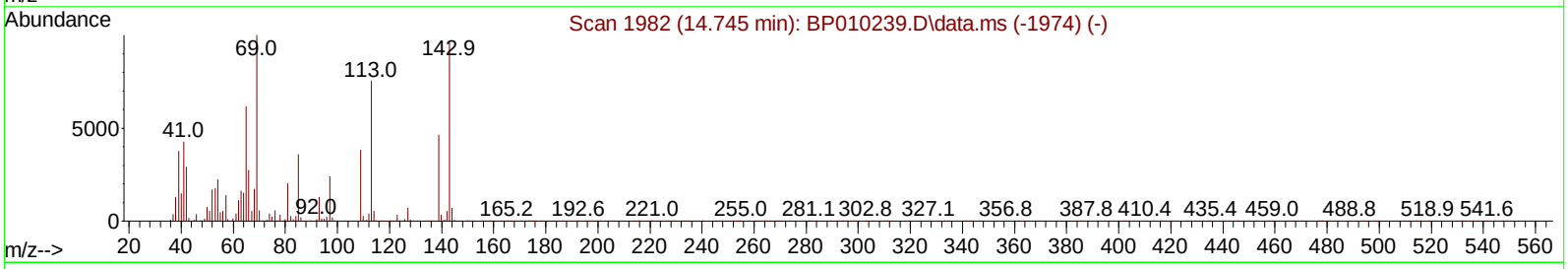

TIC: BP010244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.018) 2.30 ng/ul

response 13605

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.72#
41.00	23.20	47.44#
42.00	18.00	30.52#

Quantitation Report (Qedit)

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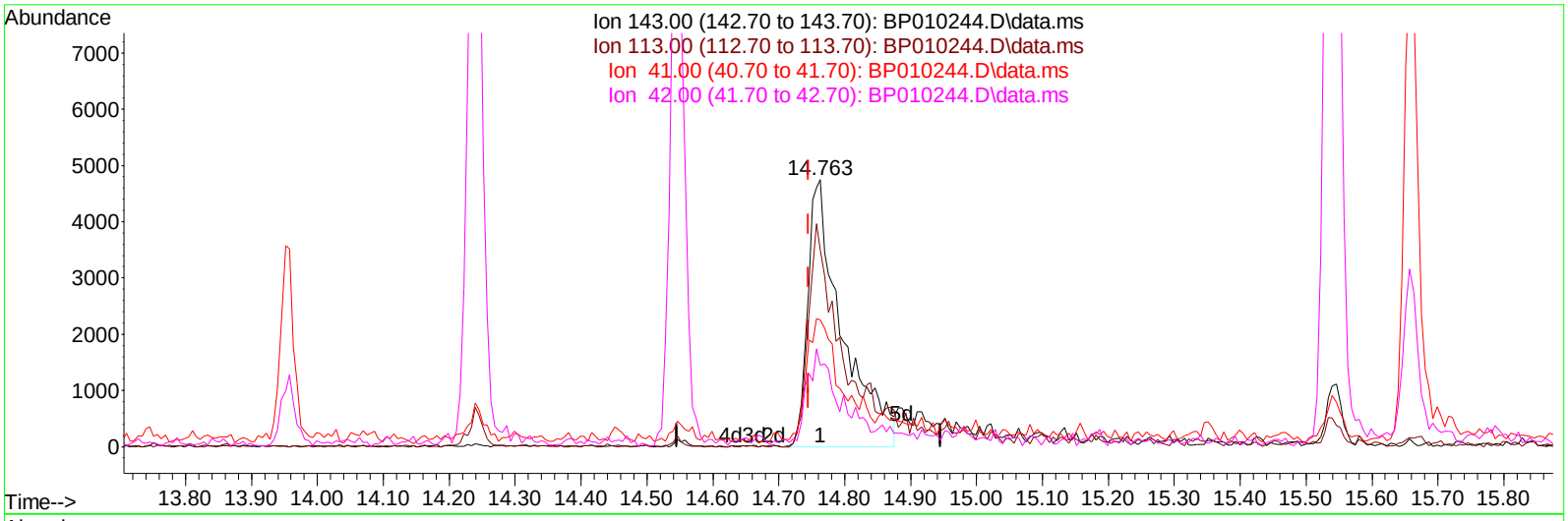

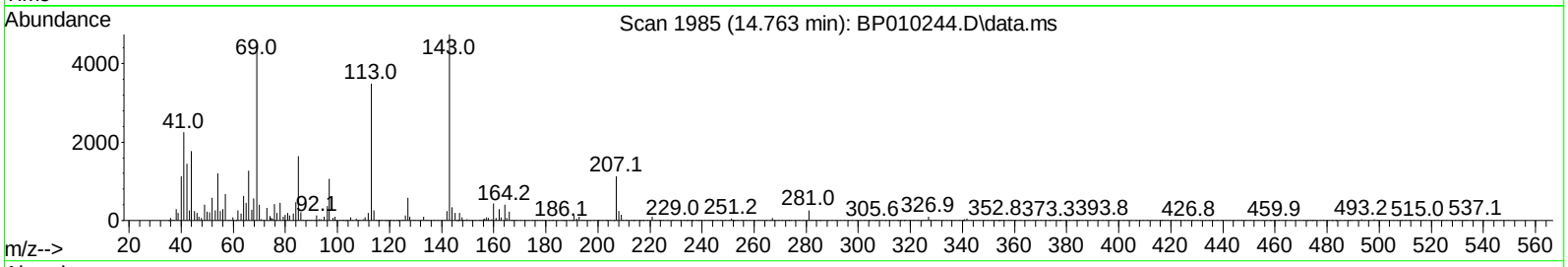

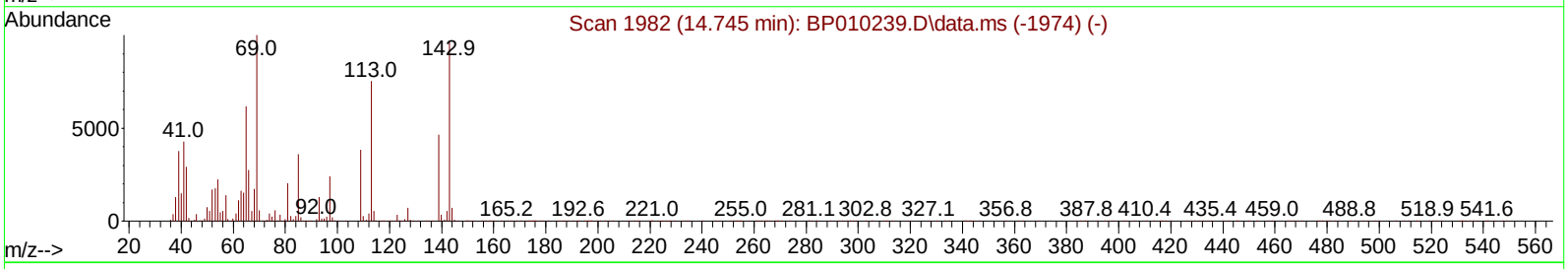

TIC: BP010244.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.018) 2.91 ng/ul m

response 17217

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.72#
41.00	23.20	47.44#
42.00	18.00	30.52#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	154271	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	653640	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.546	164	412591	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	750453	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	586585	20.000	ng/ul	# 0.00
88) Perylene-d12	23.822	264	567605	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	15962m	4.512	ng/uL	0.00
4) Pyridine-d5	3.781	84	57565	5.784	ng/ul	0.01
7) Phenol-d5	7.081	99	80294	5.722	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	317516	37.710	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	272046	26.178	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	169036	14.539	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	194659	37.444	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	197728	34.826	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	307197	27.793	ng/ul	0.00
31) 4-Chloroaniline-d4	10.858	131	351053	22.151	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1220197	38.018	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1356824	34.930	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	17217m	2.914	ng/ul	0.02
60) Fluorene-d10	15.540	176	1034280	37.463	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	135577	28.571	ng/ul	0.00
73) Anthracene-d10	17.398	188	1500823	41.338	ng/ul	0.00
81) Pyrene-d10	19.628	212	1539970	46.163	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.669	264	1297624	43.607	ng/ul	0.00

Target Compounds Qvalue

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

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