Quantitation Report (Qedit)

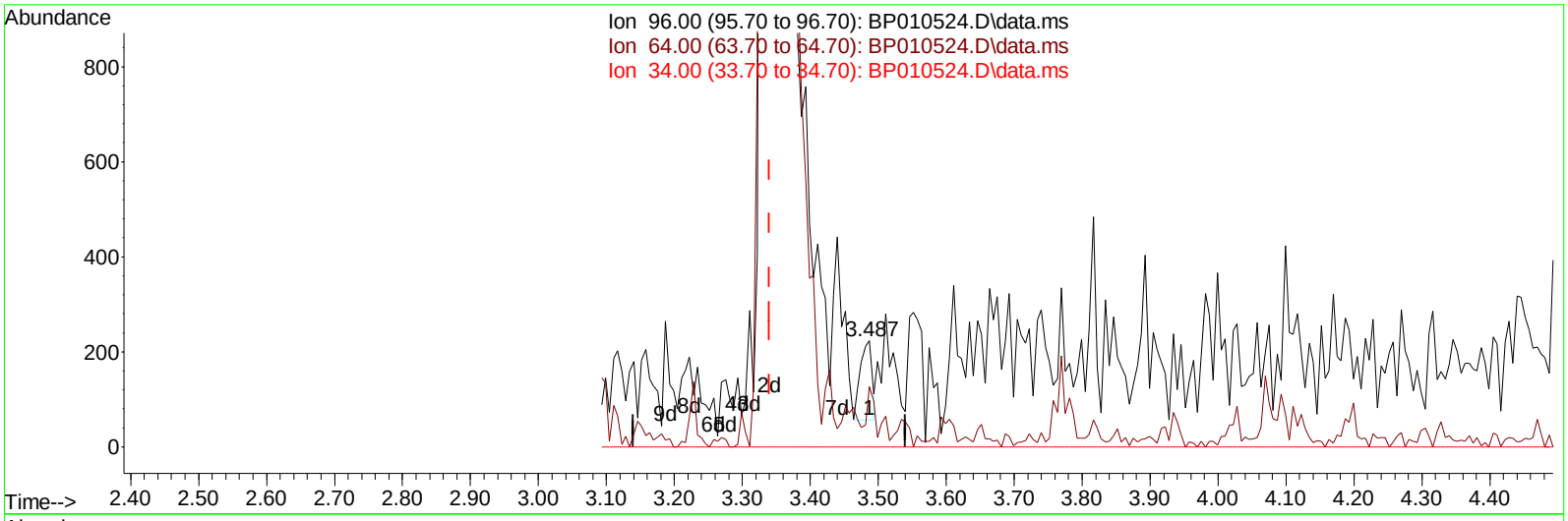
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
Data File : BP010524.D
Acq On : 24 May 2022 19:41
Operator : CG/JU
Sample : N2950-06
Misc :
ALS Vial : 53 Sample Multiplier: 1

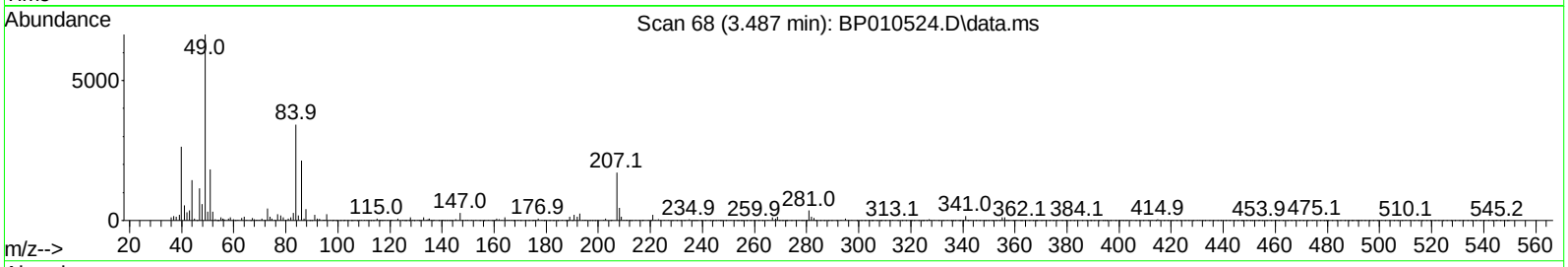
Quant Time: May 24 23:18:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 24 05:50:40 2022
Response via : Initial Calibration

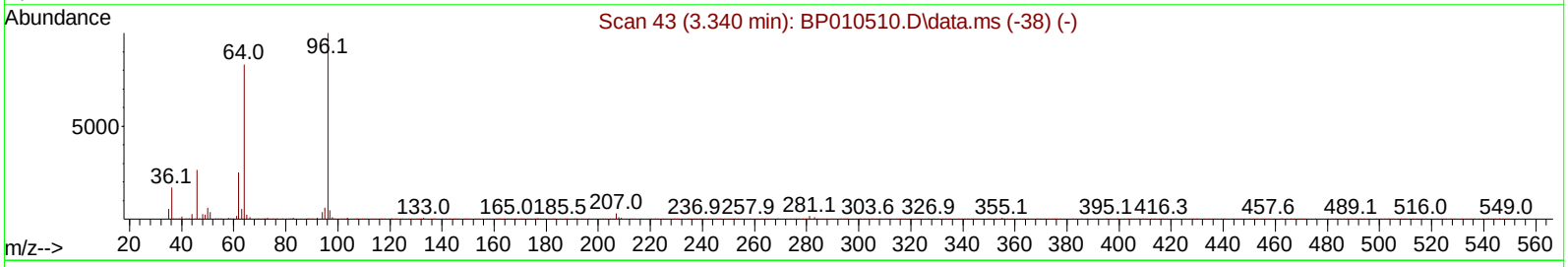
Instrument :
BNA_P
ClientSampleId :
DBTG4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/25/2022
Supervised By :Yogesh Patel 05/26/2022







TIC: BP010524.D\data.ms

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

3.487min (+ 0.147) 0.06 ng/uL

response 198

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.95
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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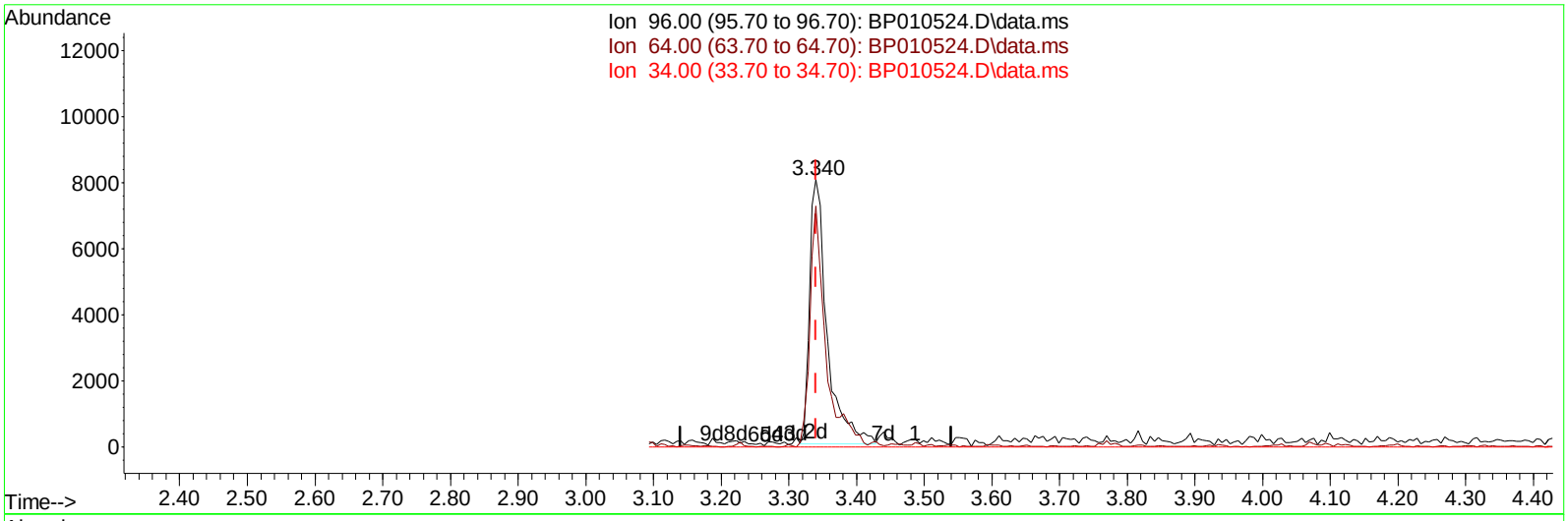

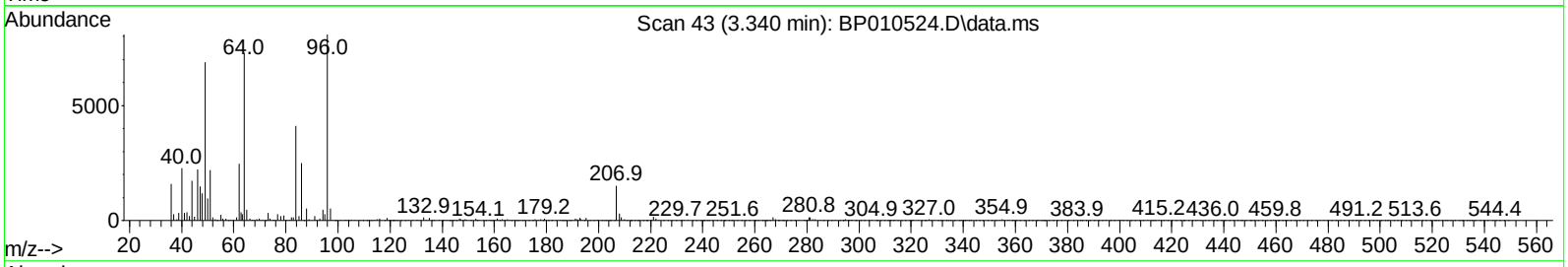

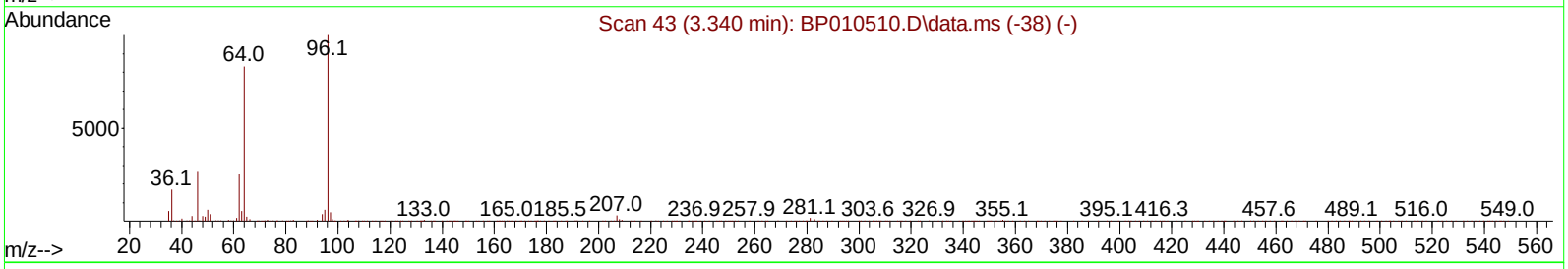

TIC: BP010524.D\data.ms

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

3.340min (-0.000) 4.43 ng/uL m

response 14487

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	90.18#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
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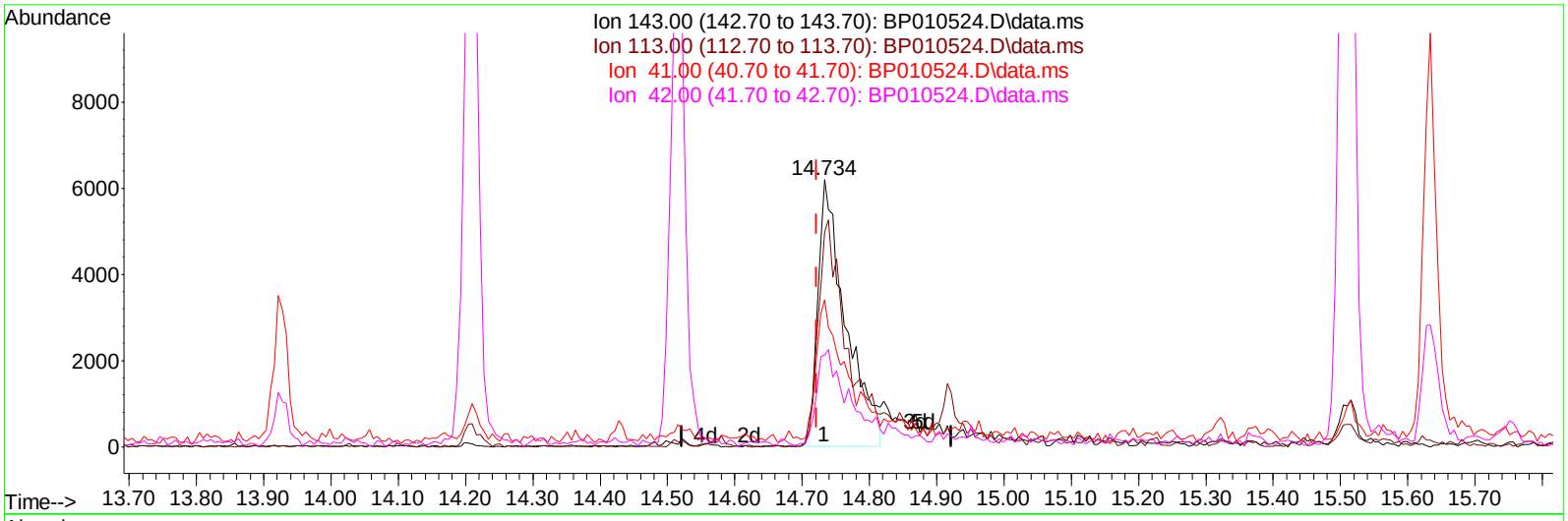

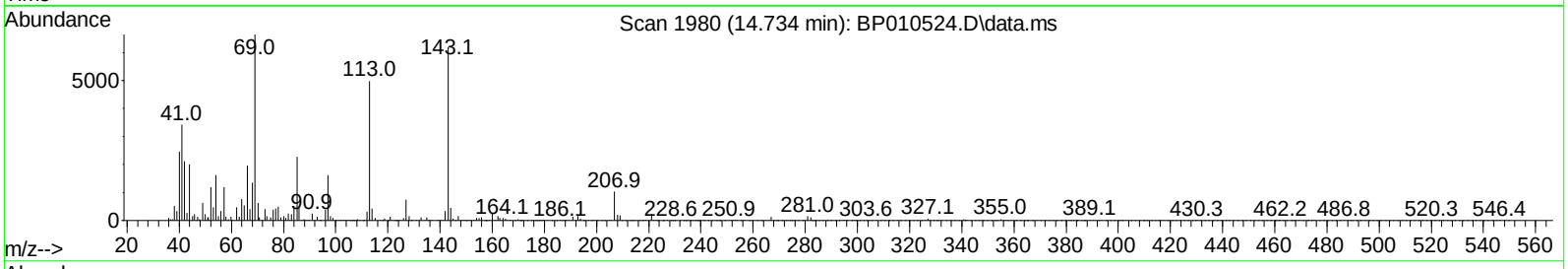

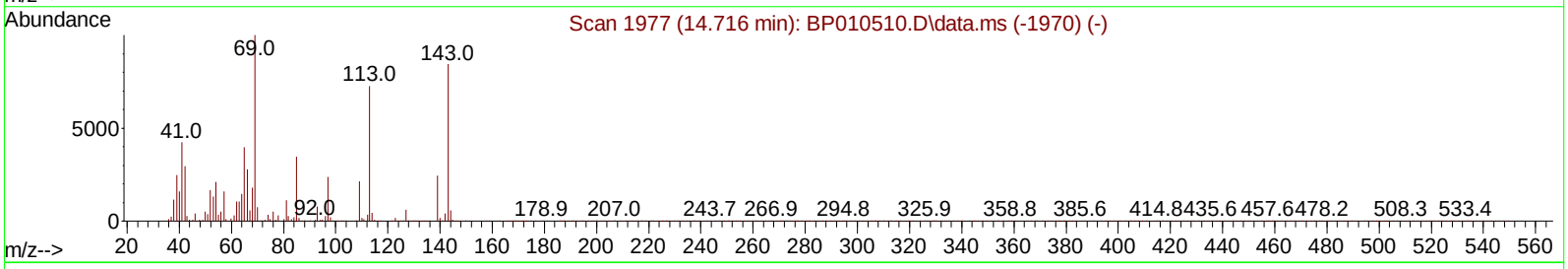

TIC: BP010524.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.012) 3.08 ng/ul

response 17727

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.46#
41.00	23.20	55.15#
42.00	18.00	34.25#

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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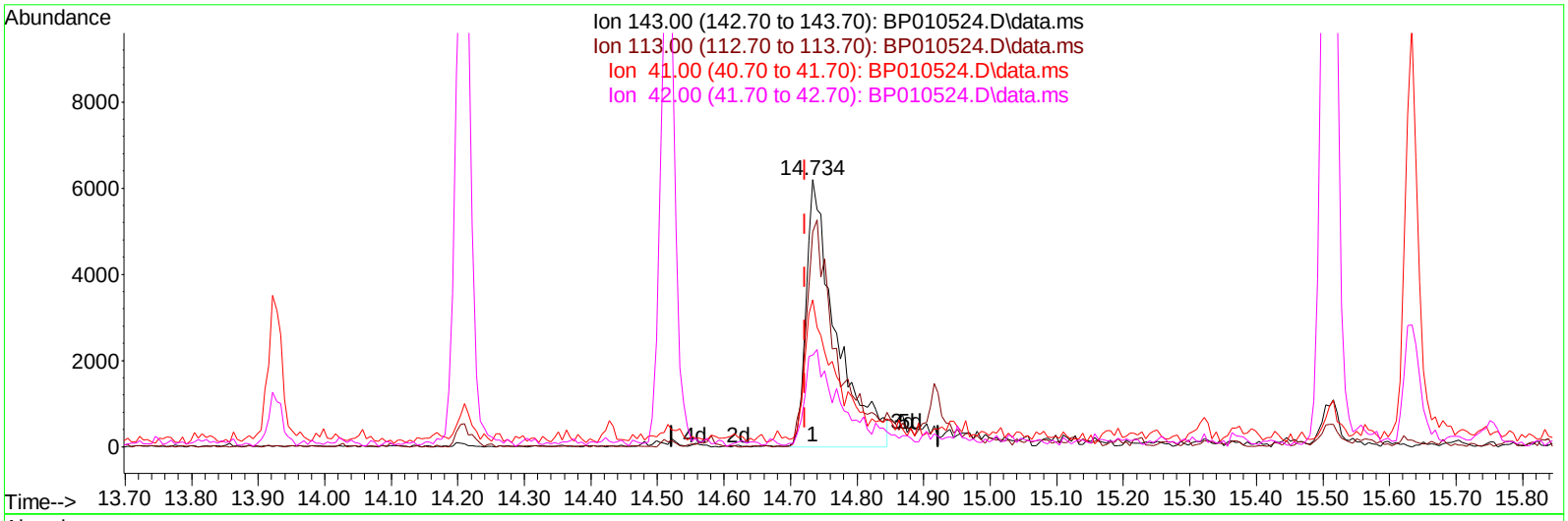

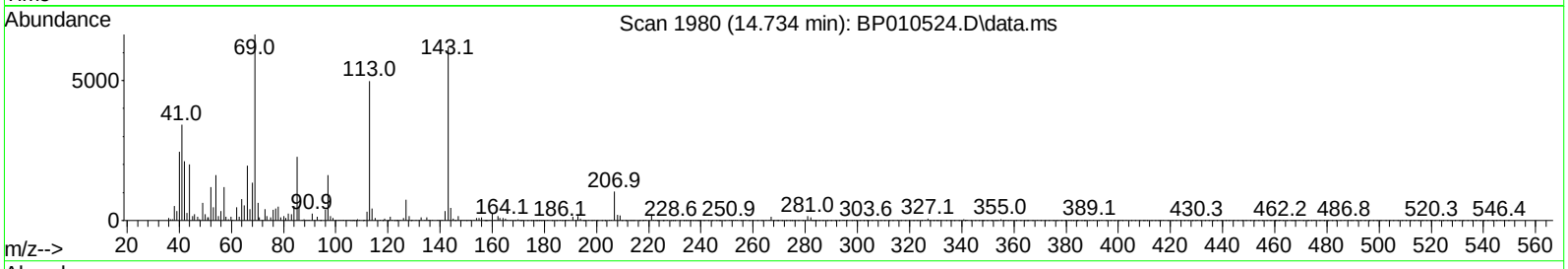

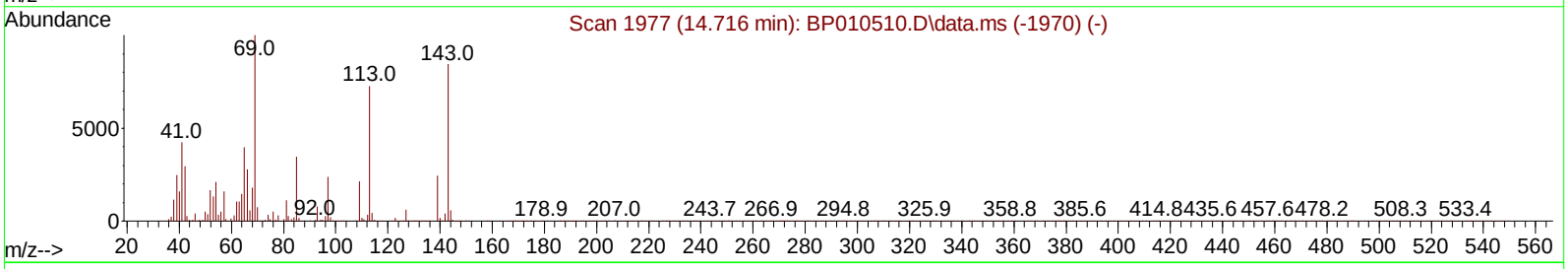

TIC: BP010524.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.012) 3.33 ng/ul m

response 19192

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.46#
41.00	23.20	55.15#
42.00	18.00	34.25#

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 05/26/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.887	152	139381	20.000	ng/ul		0.00
20) Naphthalene-d8	10.687	136	638009	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.516	164	439410	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.269	188	945217	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.351	240	769999	20.000	ng/ul		0.00
88) Perylene-d12	23.774	264	611992	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.340	96	14487m	4.430	ng/uL		0.00
4) Pyridine-d5	3.769	84	54878	5.911	ng/ul		0.01
7) Phenol-d5	7.052	99	75132	6.422	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	249210	31.707	ng/ul		0.00
11) 2-Chlorophenol-d4	7.416	132	221761	24.773	ng/ul		0.00
15) 4-Methylphenol-d8	8.593	113	157562	15.848	ng/ul		0.00
21) Nitrobenzene-d5	9.046	128	163677	33.148	ng/ul		0.00
24) 2-Nitrophenol-d4	9.769	143	168864	32.270	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.310	165	276985	27.994	ng/ul		0.00
31) 4-Chloroaniline-d4	10.822	131	356667	24.243	ng/ul		0.00
46) Dimethylphthalate-d6	13.928	166	1112138	34.352	ng/ul		0.00
49) Acenaphthylene-d8	14.210	160	1256140	33.638	ng/ul		0.00
54) 4-Nitrophenol-d4	14.734	143	19192m	3.331	ng/ul		0.01
60) Fluorene-d10	15.510	176	981669	34.854	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.633	200	118298	21.400	ng/ul		0.00
73) Anthracene-d10	17.369	188	1685406	39.378	ng/ul		0.00
81) Pyrene-d10	19.604	212	1966689	48.094	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.615	264	1245682	40.536	ng/ul		0.00

Target Compounds Qvalue

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

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