Quantitation Report (Qedit)

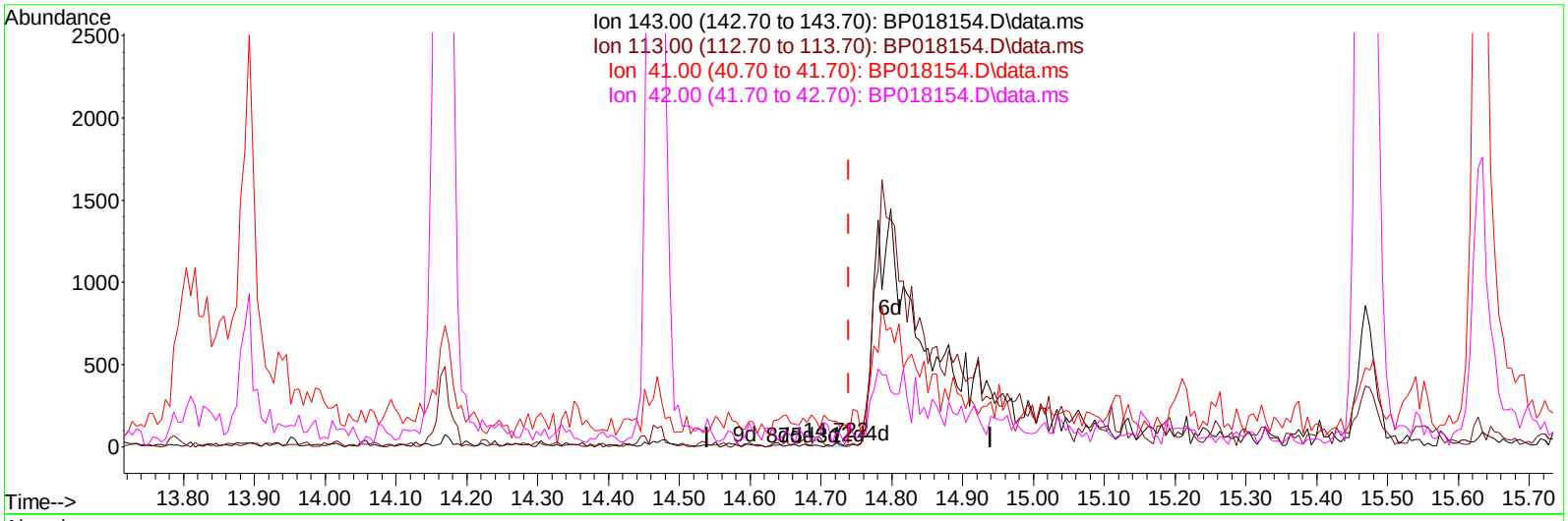
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
Data File : BP018154.D
Acq On : 14 Nov 2023 14:28
Operator : MA/JU
Sample : 05337-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

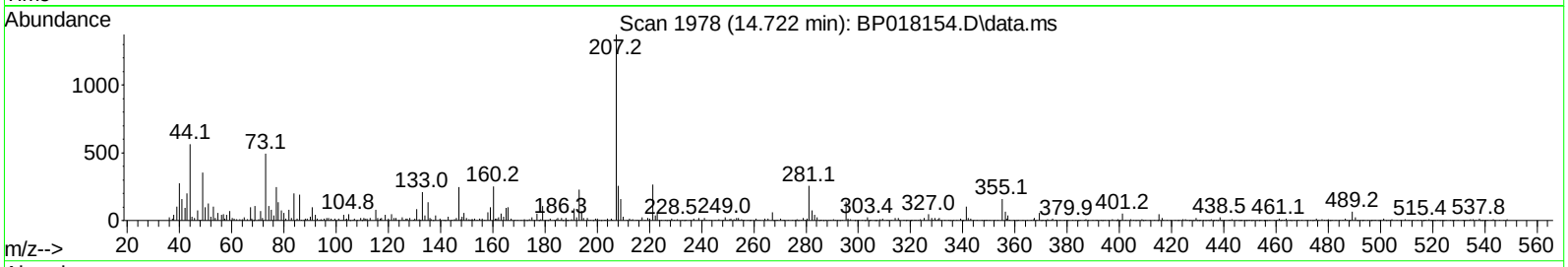
Quant Time: Nov 15 00:38:02 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 13 15:50:18 2023
Response via : Initial Calibration

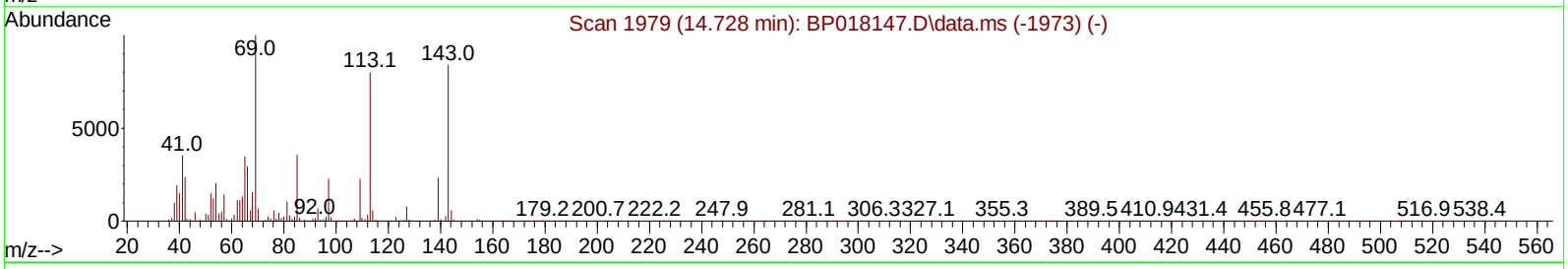
Instrument :
BNA_P
ClientSampleId :
GCDH7

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/16/2023
Supervised By :mohammad ahmed 11/16/2023







TIC: BP018154.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.018) 0.01 ng/ul

response 17

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	75.00#
41.00	47.70	560.71#
42.00	30.10	328.57#

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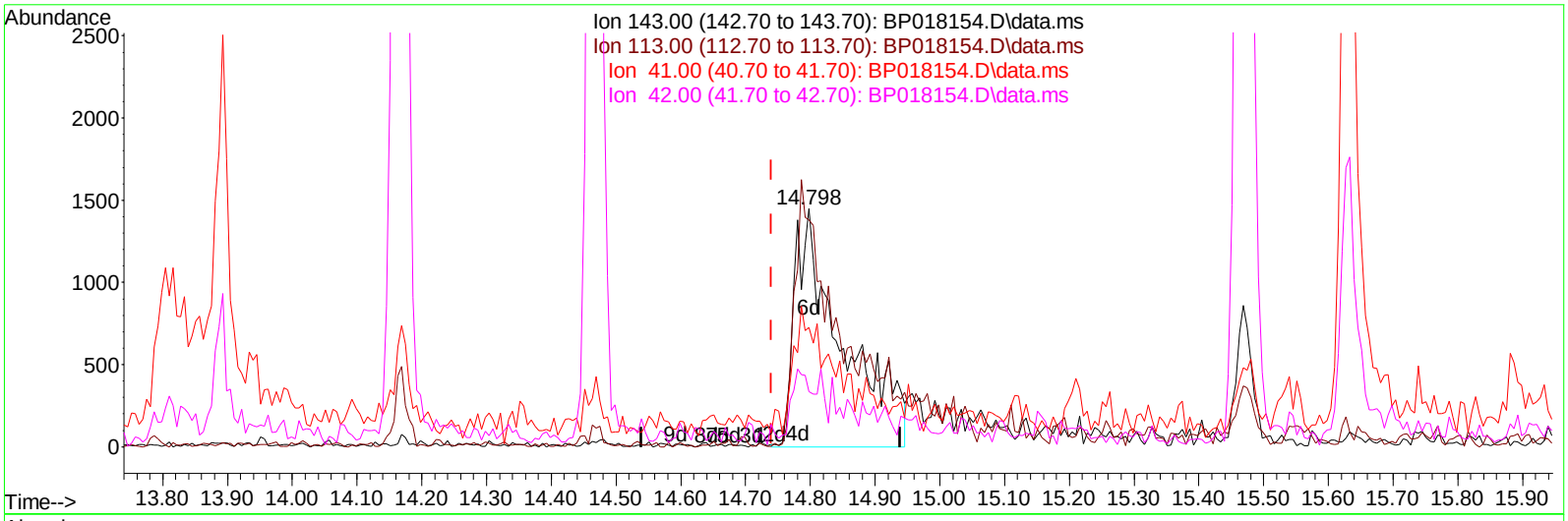

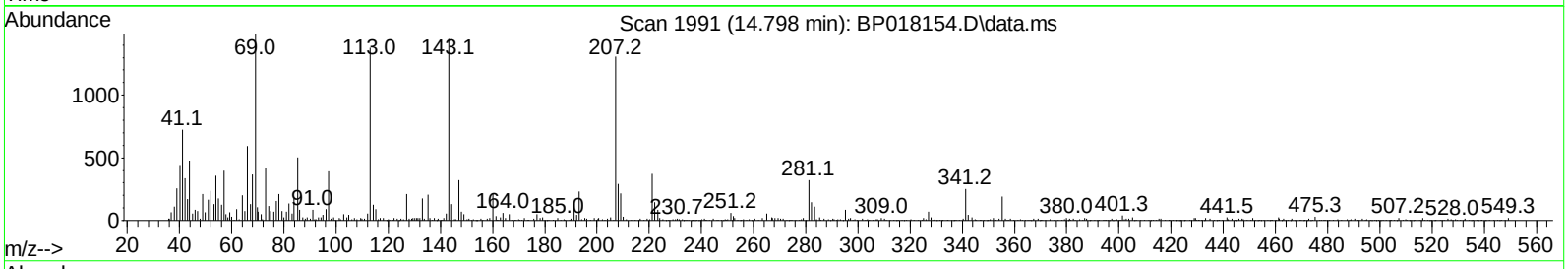

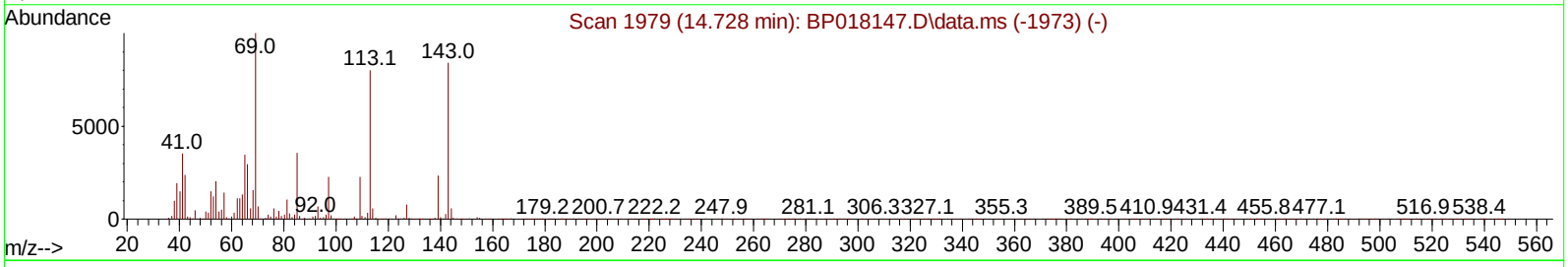

TIC: BP018154.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.059) 2.21 ng/ul m

response 7381

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	95.23
41.00	47.70	50.21
42.00	30.10	23.27#

Quantitation Report (Qedit)

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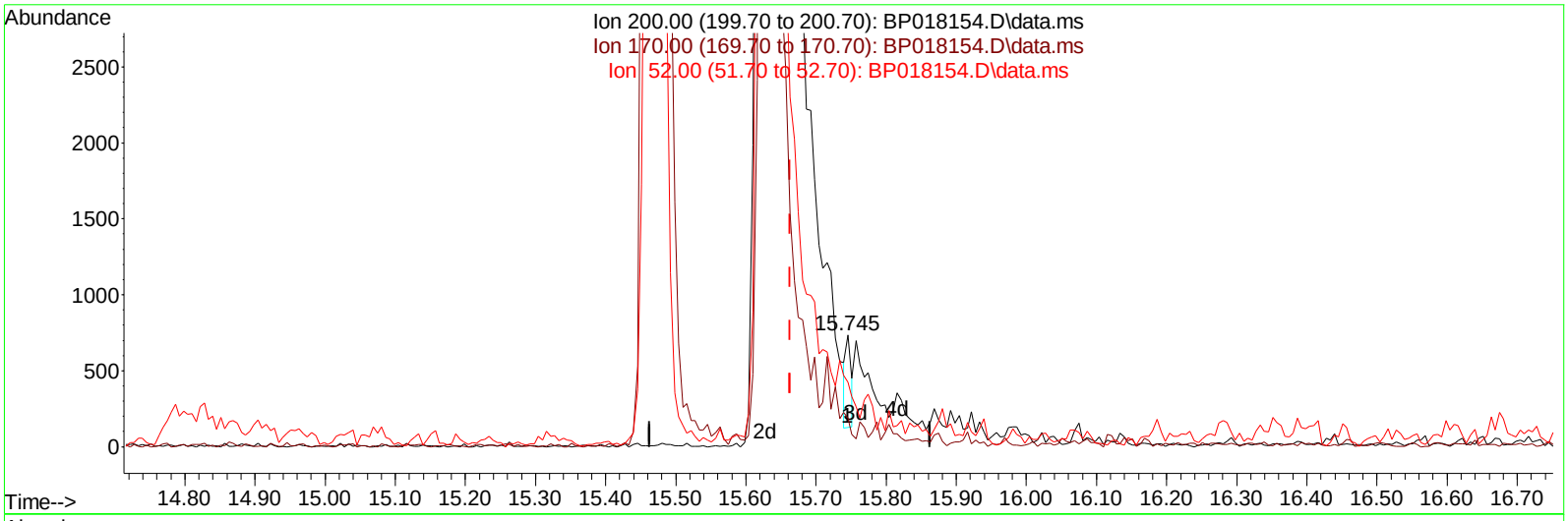

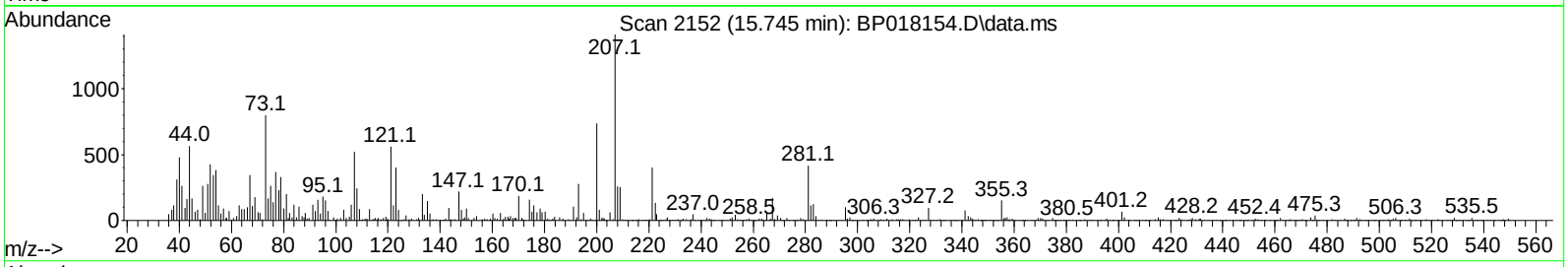

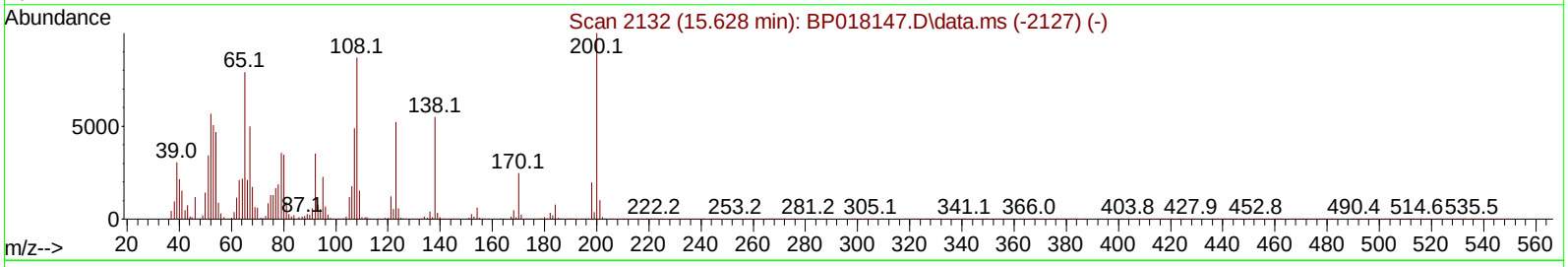

TIC: BP018154.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.745min (+ 0.082) 0.09 ng/ul

response 330

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	25.41
52.00	52.00	57.74
0.00	0.00	0.00

Quantitation Report (Qedit)

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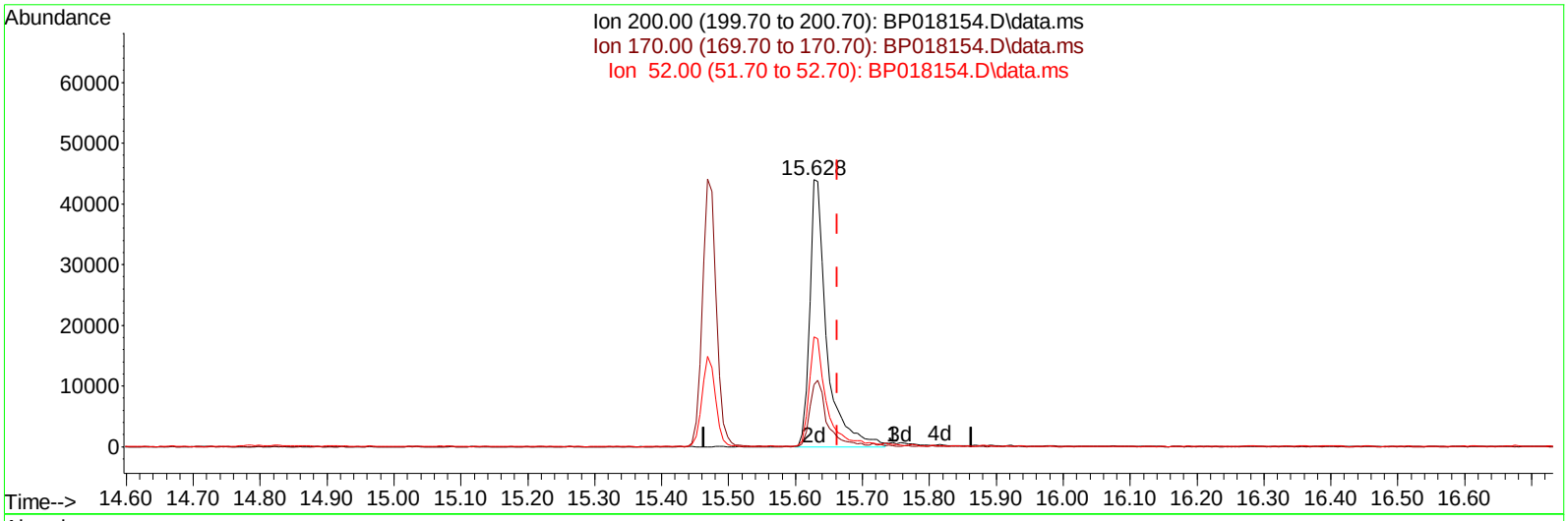

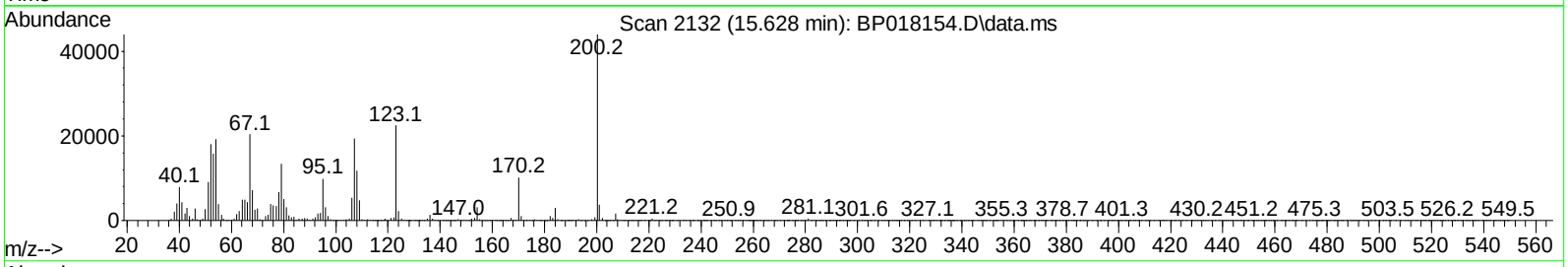

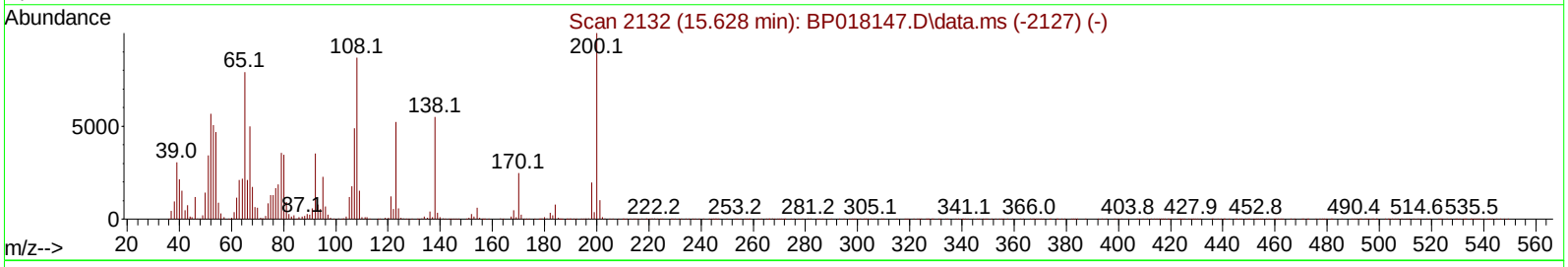

TIC: BP018154.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.628min (-0.035) 21.04 ng/ul m

response 77478

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	23.19
52.00	52.00	41.12#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	100684	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	381755	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	242318	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	509491	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	578498	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	630717	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	9120	3.195	ng/uL	0.00
4) Pyridine-d5	3.622	84	26590	3.573	ng/ul	0.02
7) Phenol-d5	6.969	99	38444	3.955	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	155263	27.180	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	138794	18.031	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	83448	10.477	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	92404	26.866	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	98217	25.206	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	153990	22.173	ng/ul	0.00
31) 4-Chloroaniline-d4	10.787	131	187506	19.688	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166	699174	31.272	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	689345	28.308	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	7381m	2.209	ng/ul	0.06
60) Fluorene-d10	15.469	176	494125	26.769	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	77478m	21.035	ng/ul	-0.04
73) Anthracene-d10	17.351	188	888150	32.232	ng/ul	0.00
81) Pyrene-d10	19.604	212	1318491	34.121	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1179288	32.156	ng/ul	-0.02

Target Compounds Qvalue

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

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