

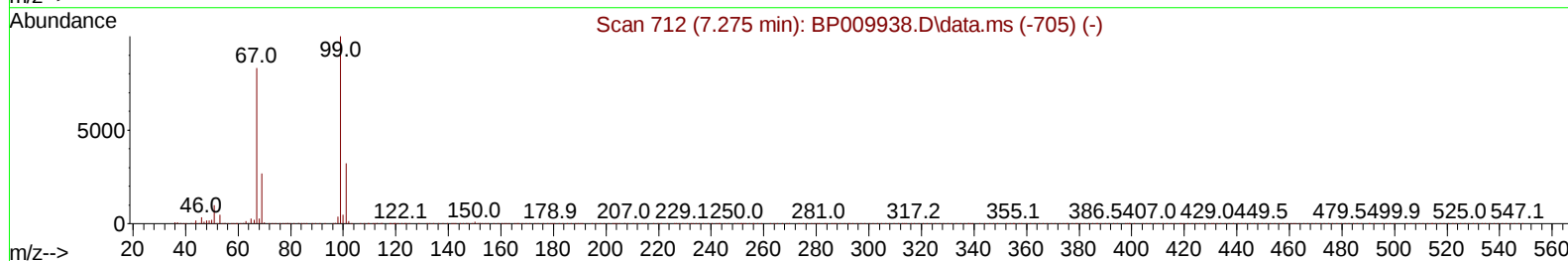
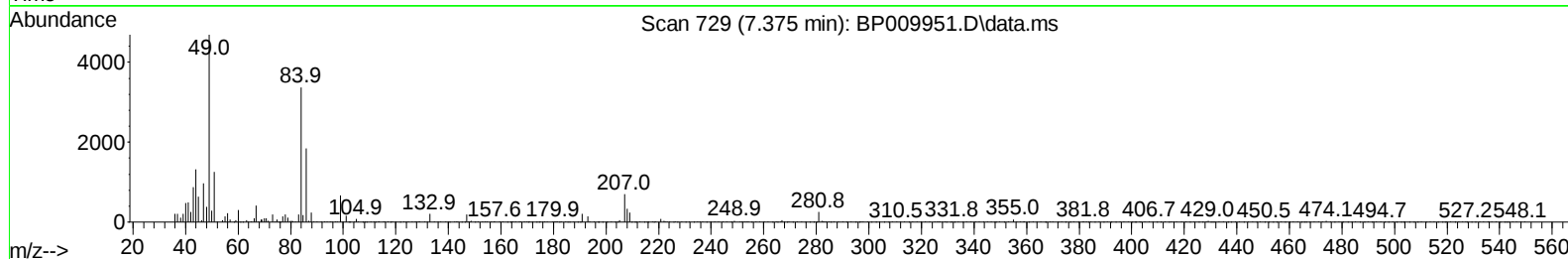
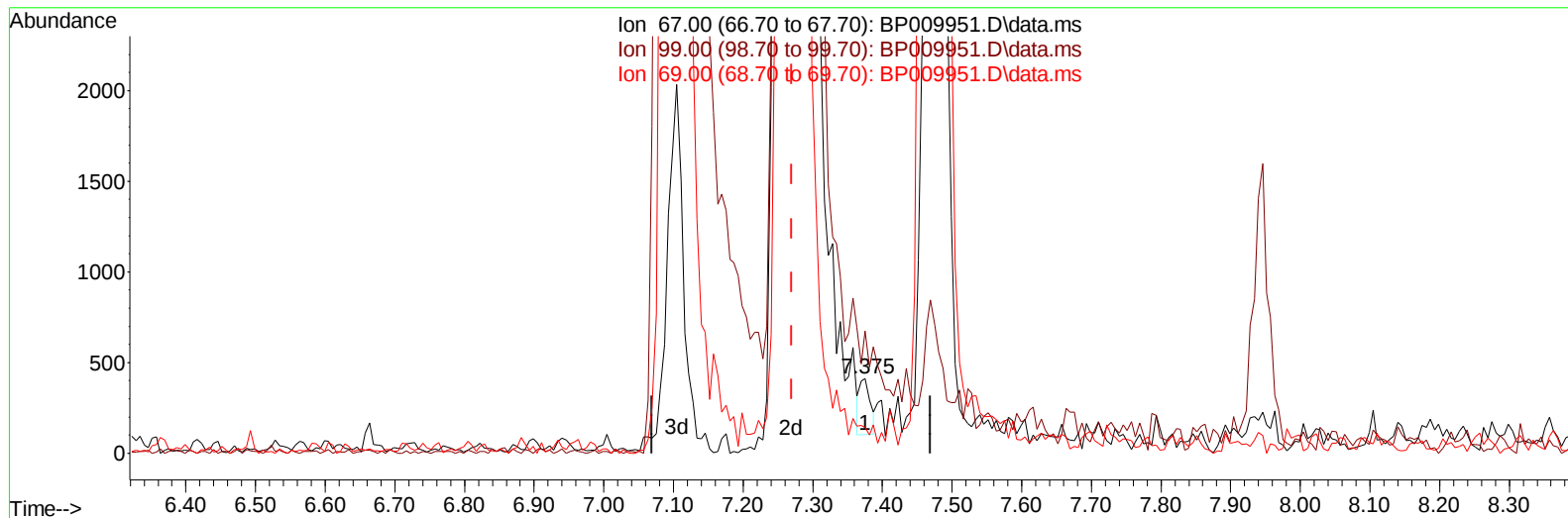
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP009951.D
 Acq On : 19 Apr 2022 18:37
 Operator : CG/JU
 Sample : PB144172BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK172

Manual IntegrationsAPPROVED

Quant Time: Apr 20 01:32:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/20/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP009951.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.375min (+ 0.106) 0.04 ng/u1

response 325

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	163.59
69.00	40.10	34.71
0.00	0.00	0.00

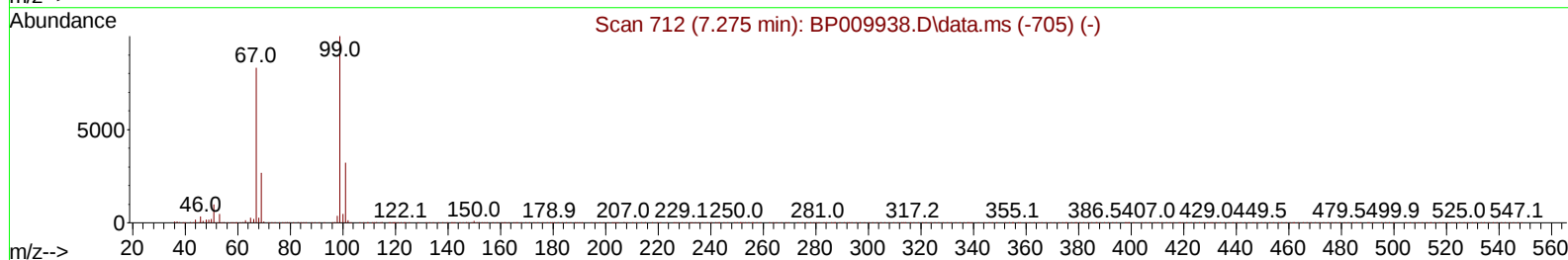
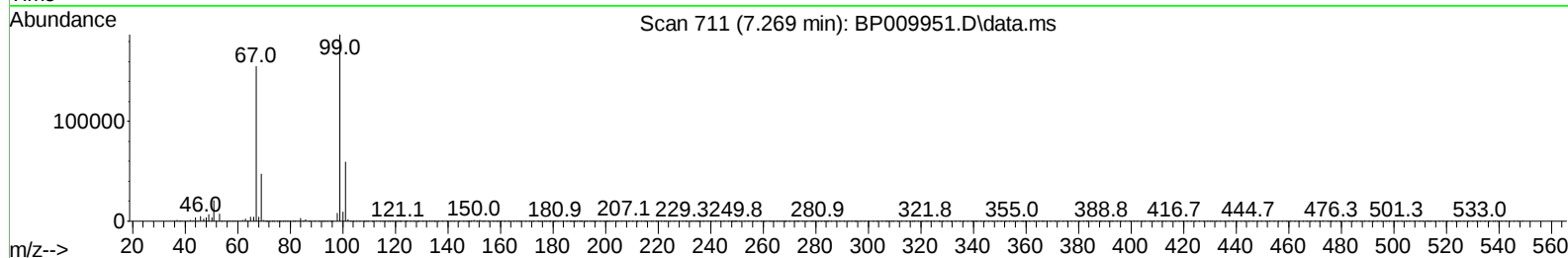
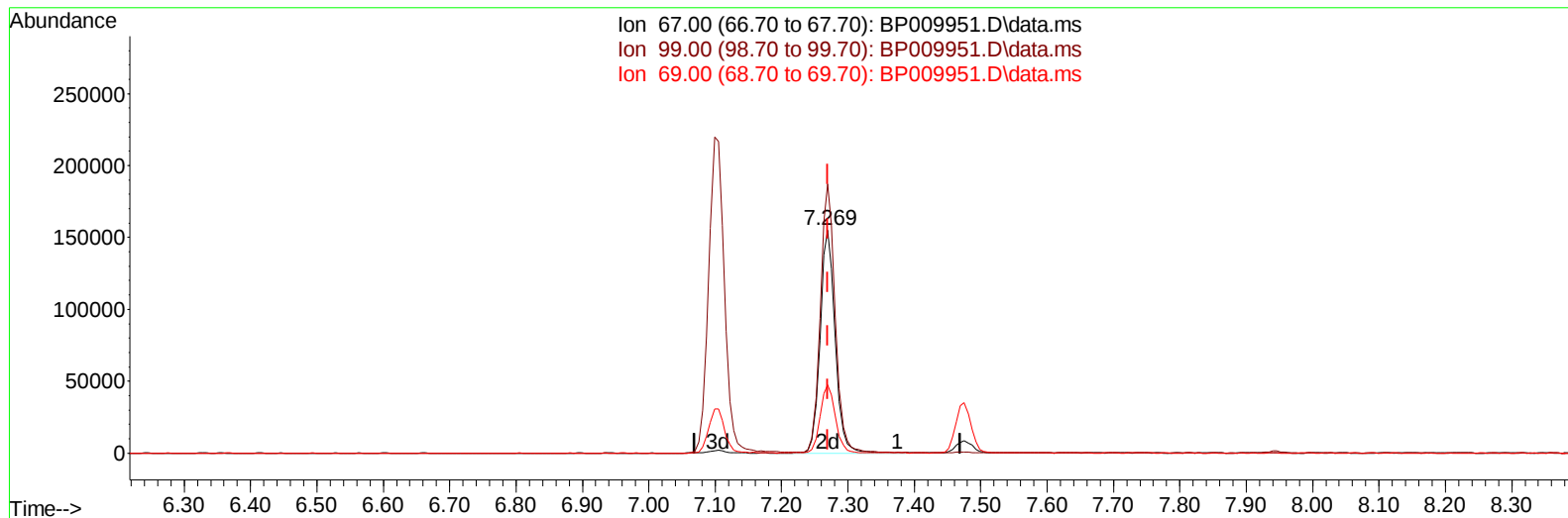
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TIC: BP009951.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.269min (-0.000) 31.10 ng/u1 m

response 245834

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	120.59#
69.00	40.10	30.68#
0.00	0.00	0.00

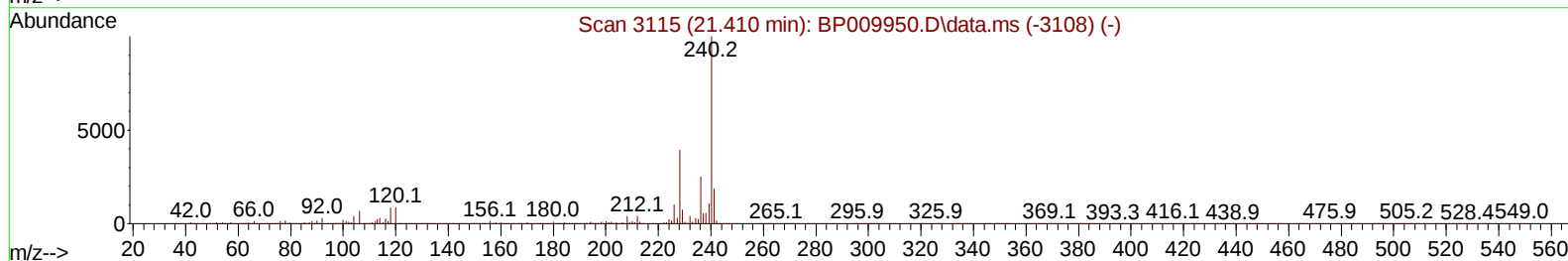
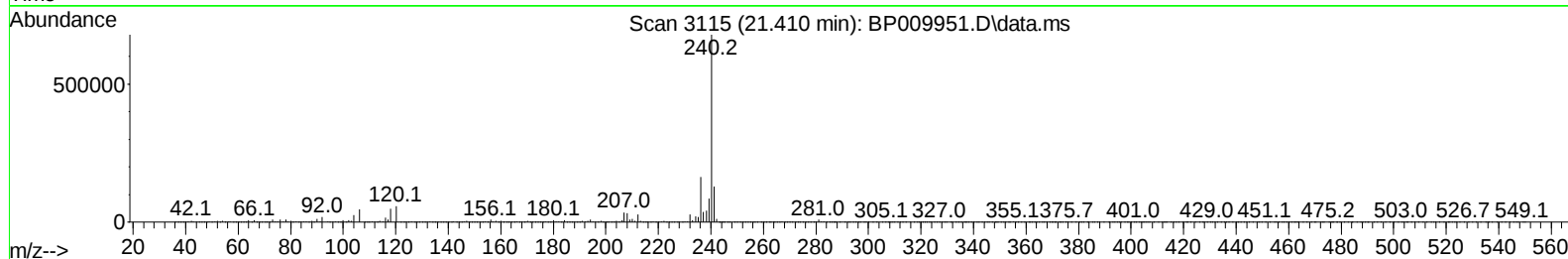
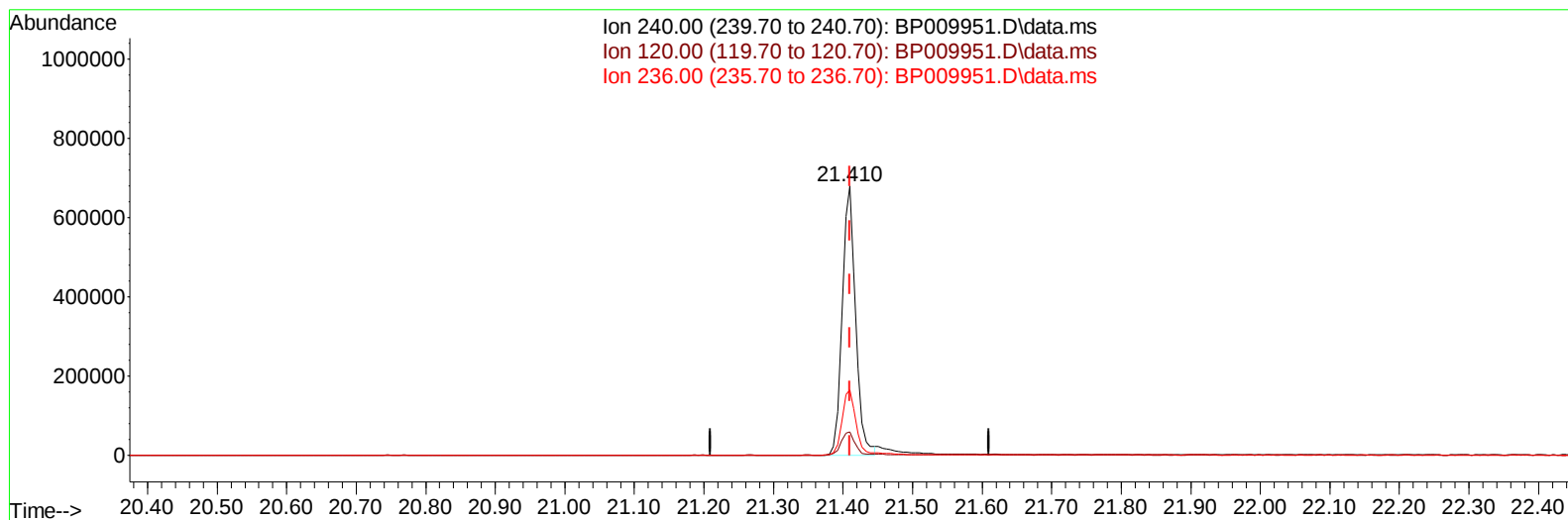
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ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP009951.D\data.ms

(79) Chrysene-d12 (I)

21.410min (-0.000) 20.00 ng/u1

response 911648

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.54
236.00	31.60	24.21#
0.00	0.00	0.00

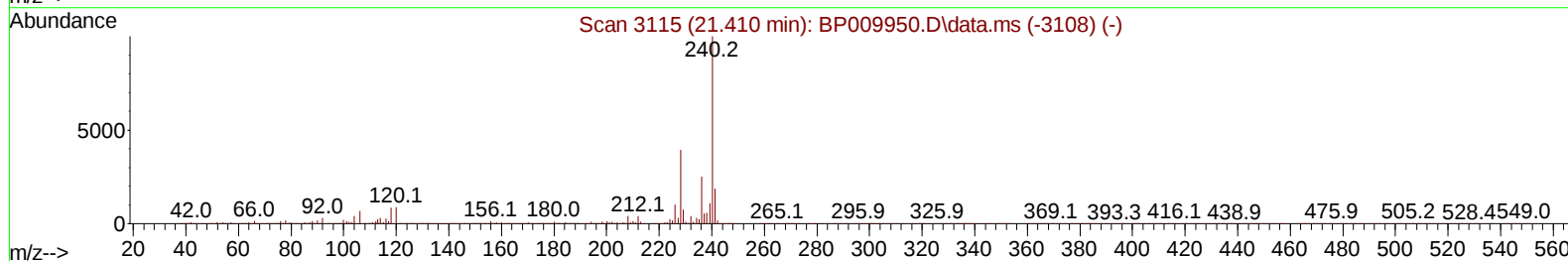
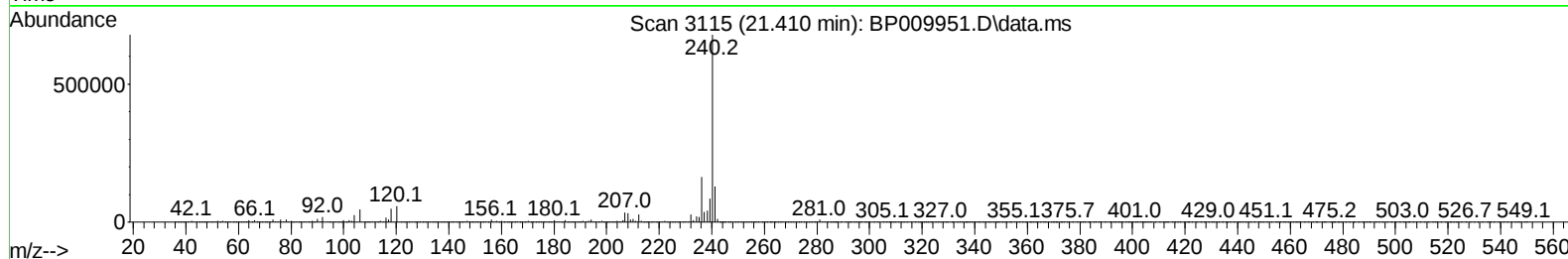
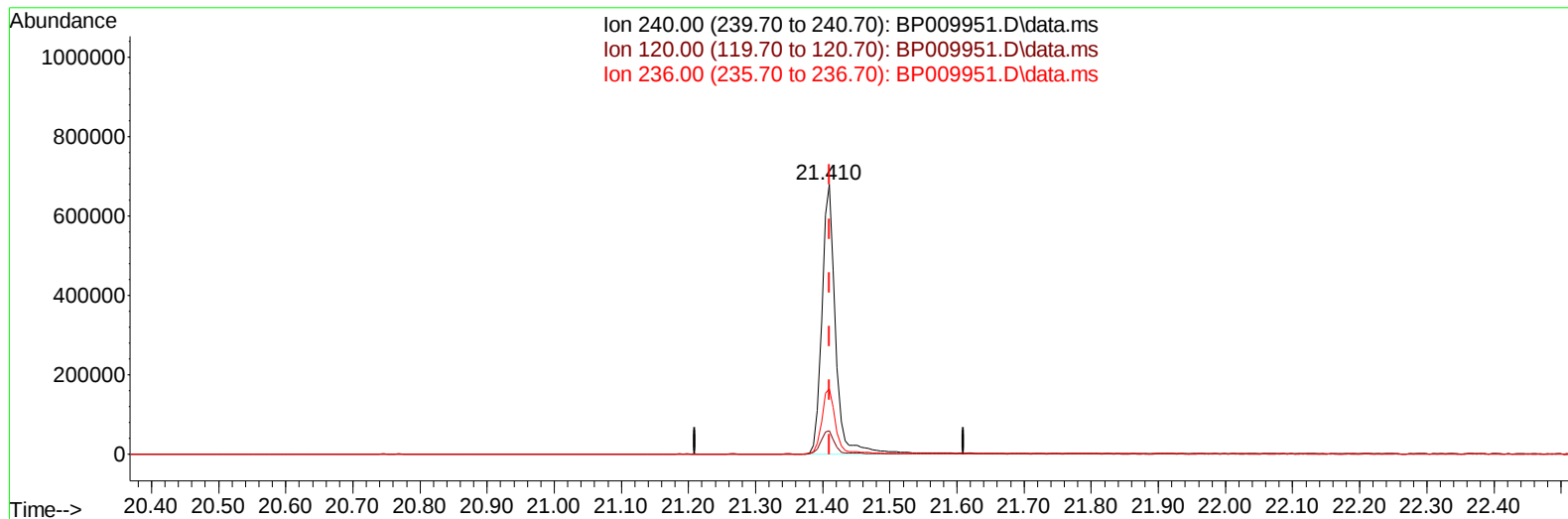
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
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TIC: BP009951.D\data.ms

(79) Chrysene-d12 (I)

21.410min (-0.000) 20.00 ng/ul m

response 954936

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.54
236.00	31.60	24.21#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	151246	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	649751	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	459198	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.328	188	1000645	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	954936m	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	858955	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	21160	6.244	ng/uL	0.00
4) Pyridine-d5	3.787	84	266019	28.470	ng/ul	0.00
7) Phenol-d5	7.099	99	362764	27.320	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	245834m	31.098	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	303333	30.284	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	312485	28.191	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	156078	33.311	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	176165	34.293	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	302585	28.149	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	440195	28.752	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1151556	31.925	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1241681	29.056	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	169360	25.890	ng/ul	0.00
60) Fluorene-d10	15.569	176	950678	31.330	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	167652	28.282	ng/ul	0.00
73) Anthracene-d10	17.428	188	1529051	31.915	ng/ul	0.00
81) Pyrene-d10	19.657	212	1853144	35.532	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1451395	32.447	ng/ul	0.00

Target Compounds Qvalue

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

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