

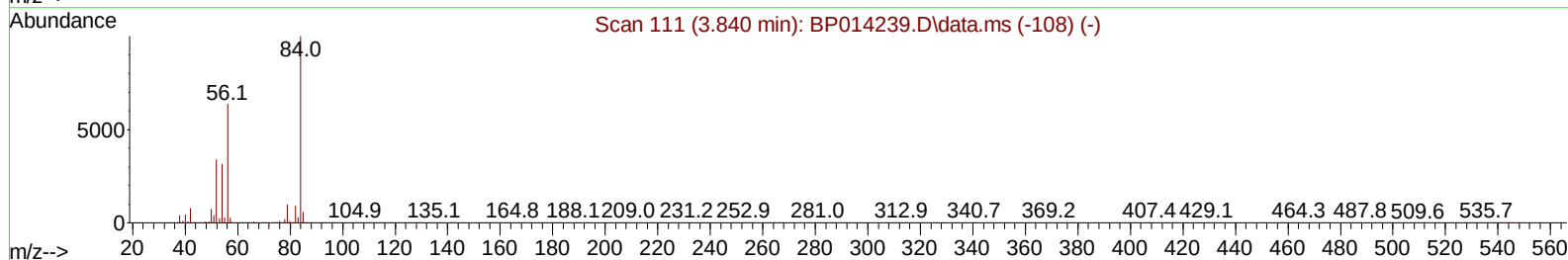
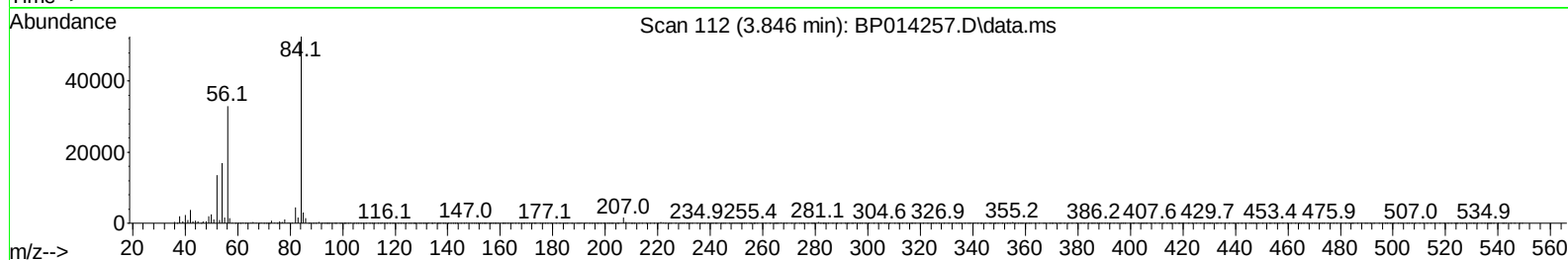
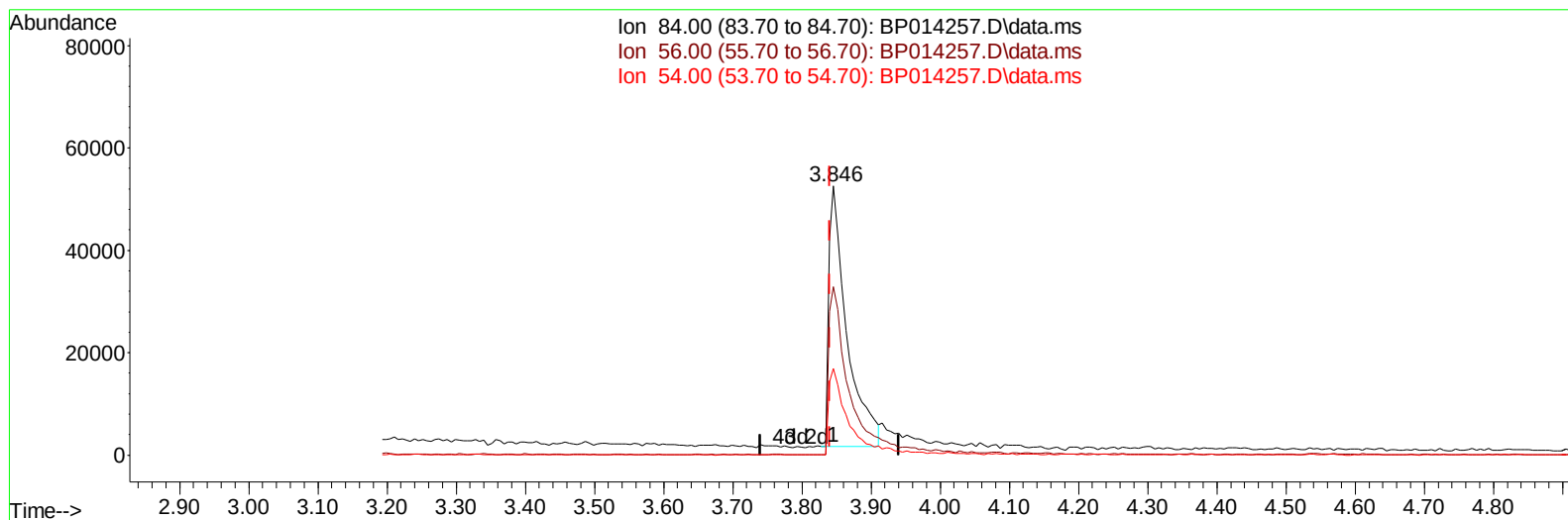
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014257.D
 Acq On : 23 Mar 2023 14:42
 Operator : CG/JU
 Sample : 02010-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB944

Manual IntegrationsAPPROVED

Quant Time: Mar 24 02:05:31 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 23 04:56:06 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/24/2023
 Supervised By :Jagrut Upadhyay 03/24/2023



TIC: BP014257.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 8.37 ng/ul

response 91606

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	62.72
54.00	33.20	32.22
0.00	0.00	0.00

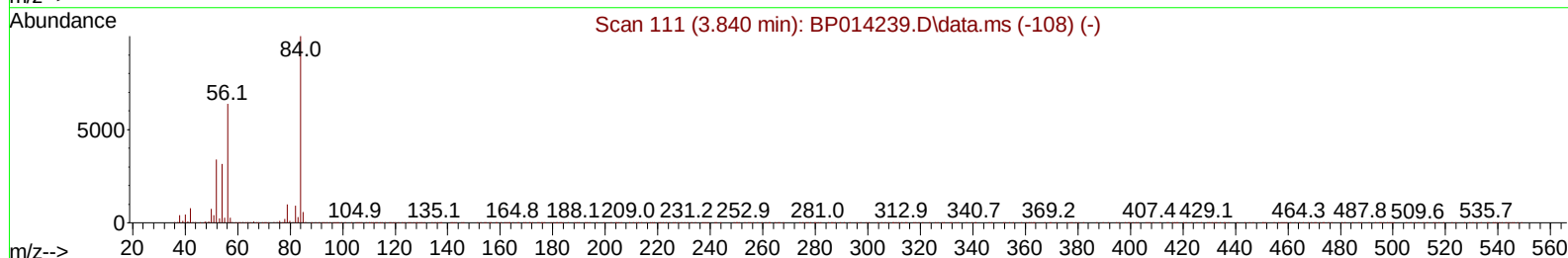
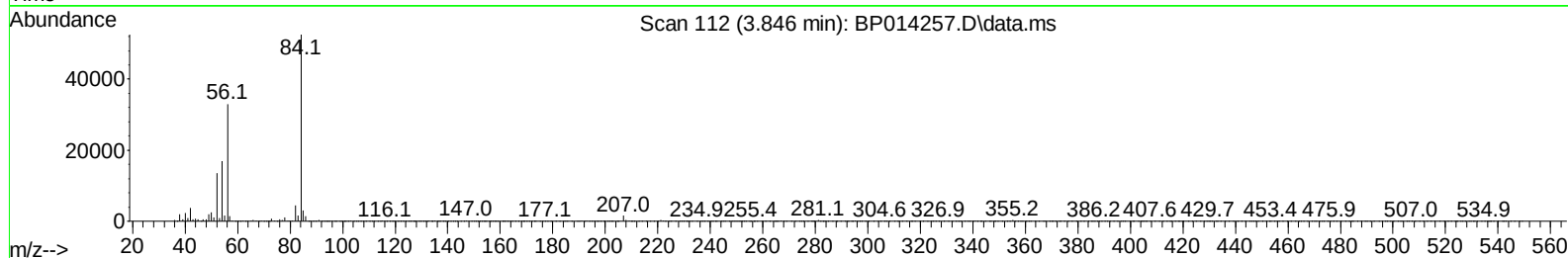
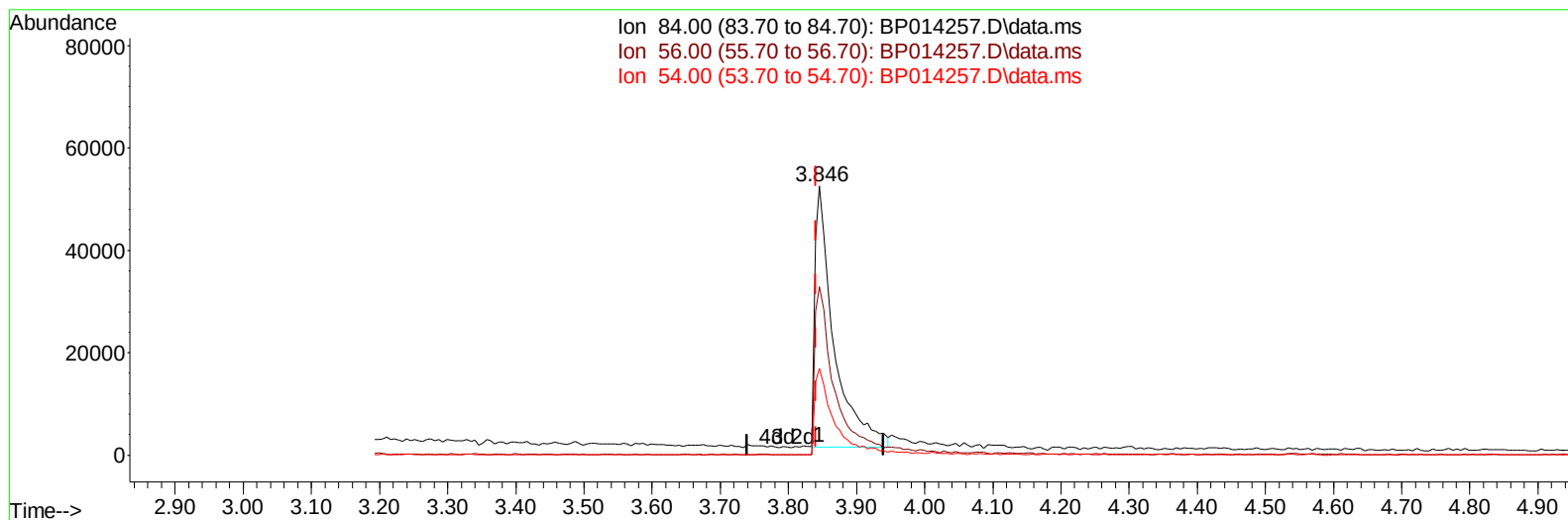
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(4) Pyridine-d5 (S)

3.846min (+ 0.006) 9.03 ng/ul m

response 98873

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	62.72
54.00	33.20	32.22
0.00	0.00	0.00

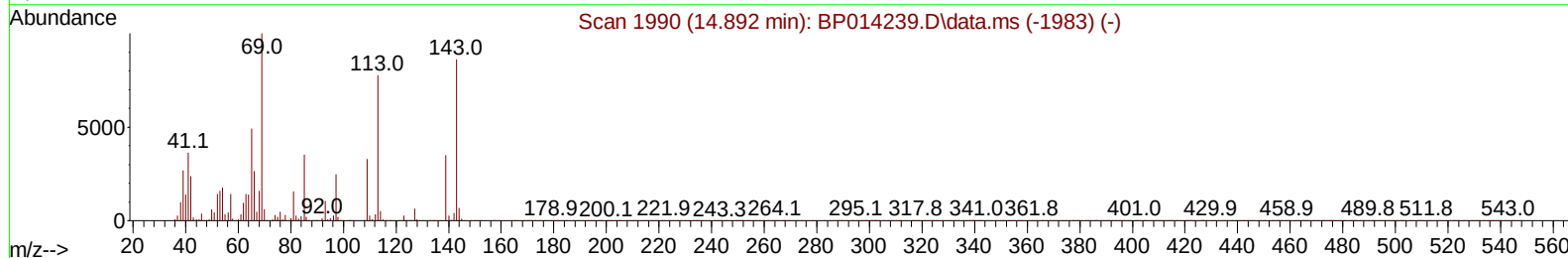
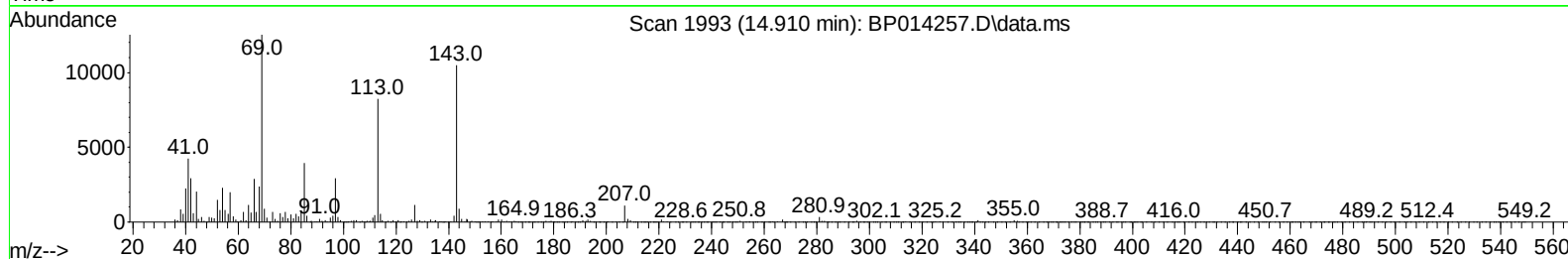
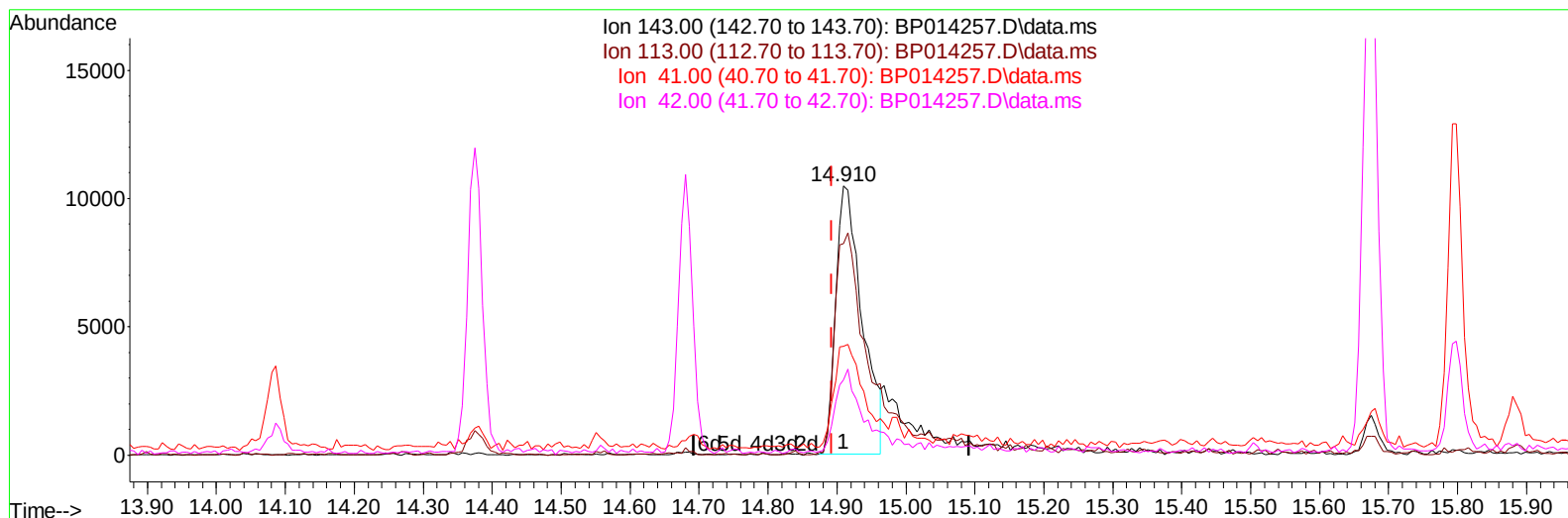
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TIC: BP014257.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.017) 3.80 ng/ul

response 27844

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	78.70#
41.00	40.20	40.46
42.00	28.50	28.07

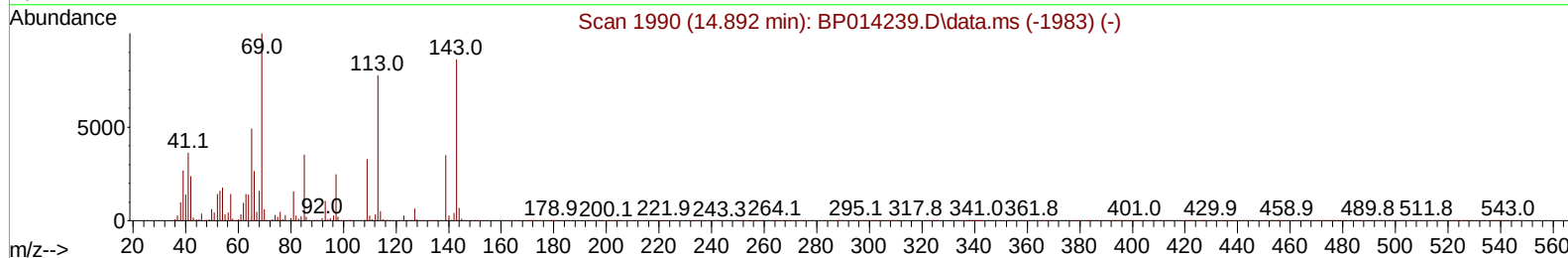
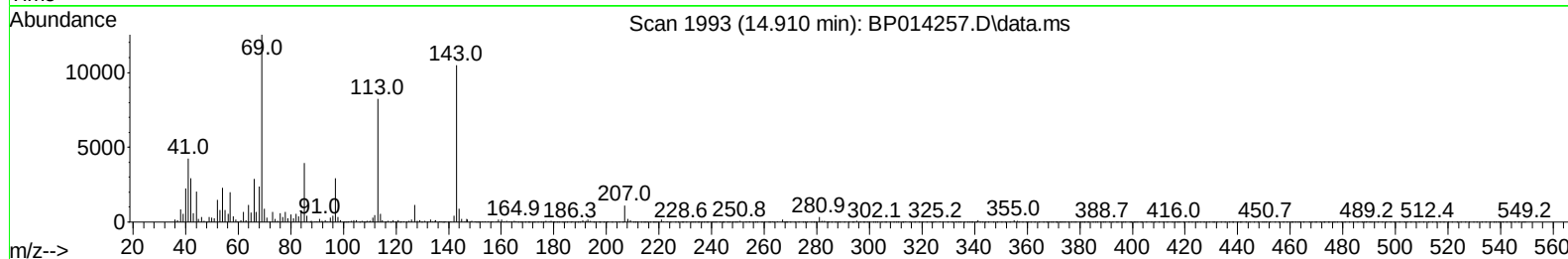
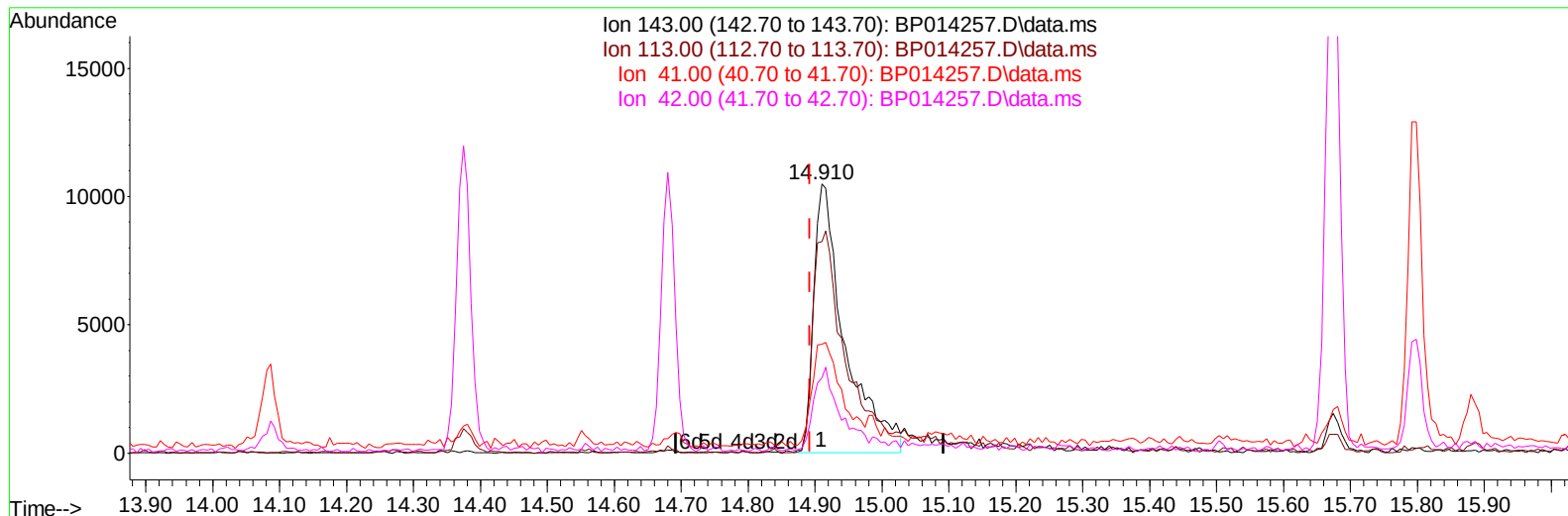
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Operator : CG/JU
Sample : 02010-01
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP014257.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.017) 4.62 ng/ul m

response 33807

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	78.70#
41.00	40.20	40.46
42.00	28.50	28.07

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	149431	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	678114	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.680	164	489308	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1144740	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.521	240	1121849	20.000	ng/ul	# 0.00
88) Perylene-d12	24.039	264	1135473	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.404	96	18271	4.620	ng/uL	0.00
4) Pyridine-d5	3.846	84	98873m	9.029	ng/ul	0.00
7) Phenol-d5	7.198	99	84310	6.393	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.357	67	258896	33.664	ng/ul	0.00
11) 2-Chlorophenol-d4	7.563	132	227719	22.666	ng/ul	0.00
15) 4-Methylphenol-d8	8.751	113	174256	16.177	ng/ul	0.00
21) Nitrobenzene-d5	9.210	128	158598	29.312	ng/ul	0.00
24) 2-Nitrophenol-d4	9.934	143	180391	27.852	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	318909	26.169	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	386685	23.977	ng/ul	0.00
46) Dimethylphthalate-d6	14.086	166	1367193	32.923	ng/ul	0.00
49) Acenaphthylene-d8	14.374	160	1399189	31.807	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	33807m	4.620	ng/ul	0.02
60) Fluorene-d10	15.674	176	1159618	32.950	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	232828	26.910	ng/ul	0.00
73) Anthracene-d10	17.539	188	1910382	34.377	ng/ul	0.00
81) Pyrene-d10	19.762	212	2306854	37.283	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.880	264	2136978	35.381	ng/ul	0.00

Target Compounds Qvalue

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

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