

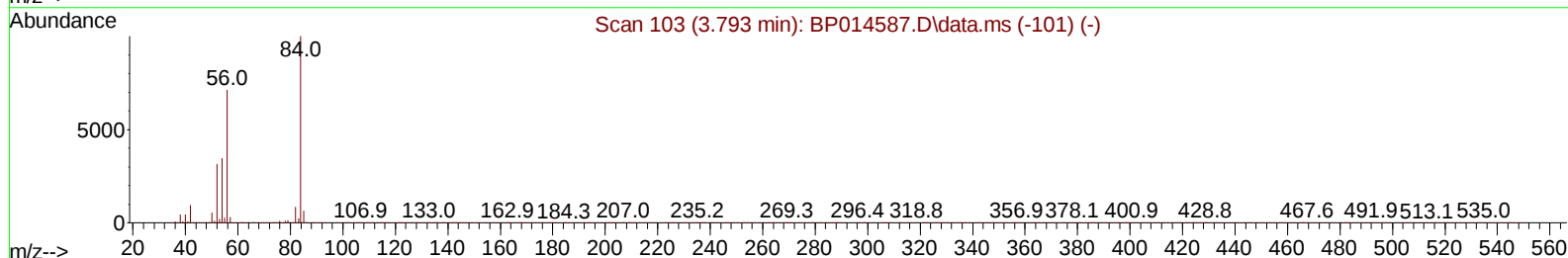
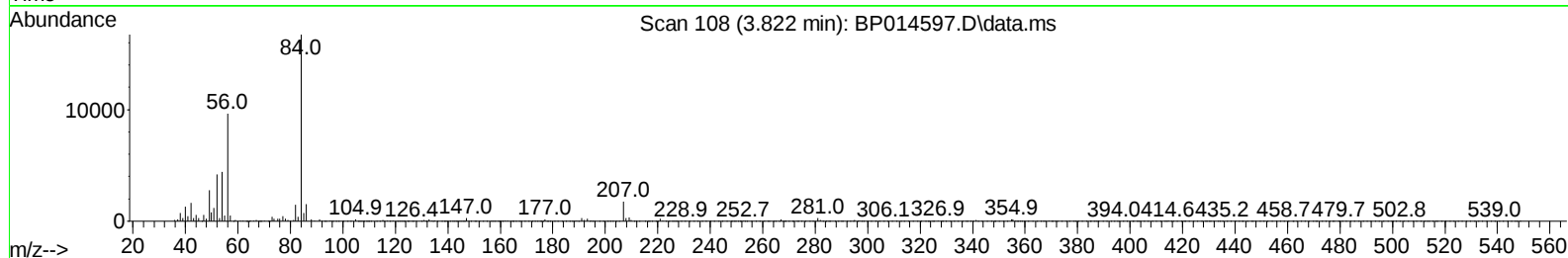
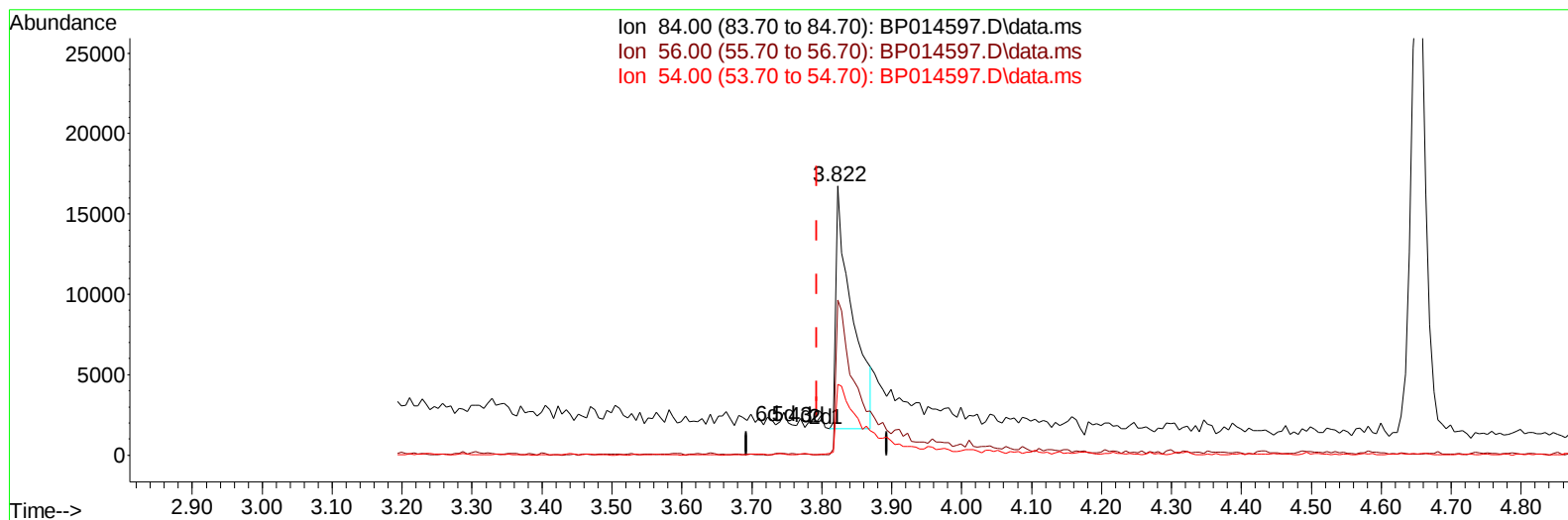
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014597.D
Acq On : 06 Apr 2023 20:39
Operator : CG/JU
Sample : 02194-11
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHDW8

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/07/2023
Supervised By :Jagrut Upadhyay 04/07/2023

Quant Time: Apr 07 00:47:45 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 07 00:45:34 2023
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TIC: BP014597.D\data.ms

(4) Pyridine-d5 (S)

3.822min (+ 0.029) 2.57 ng/u1

response 24254

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	57.76
54.00	32.60	26.40
0.00	0.00	0.00

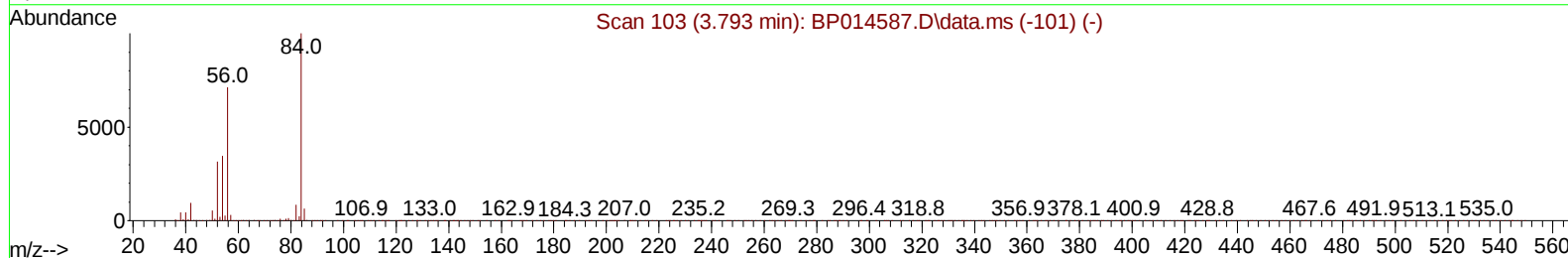
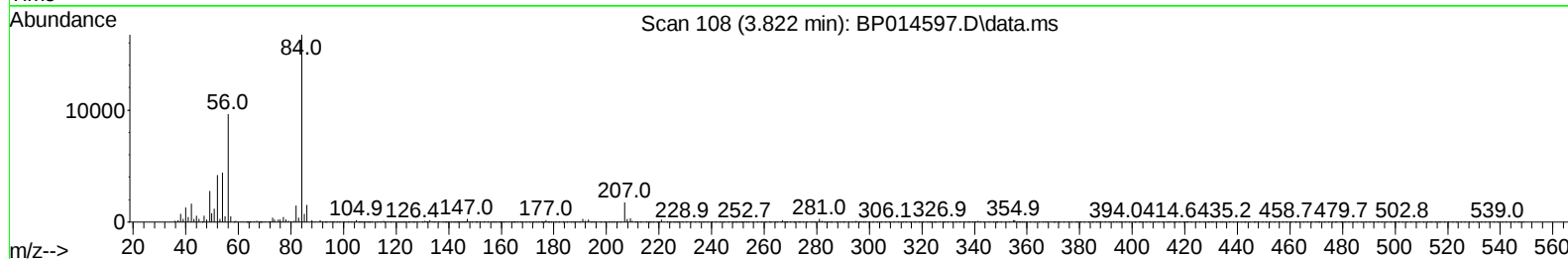
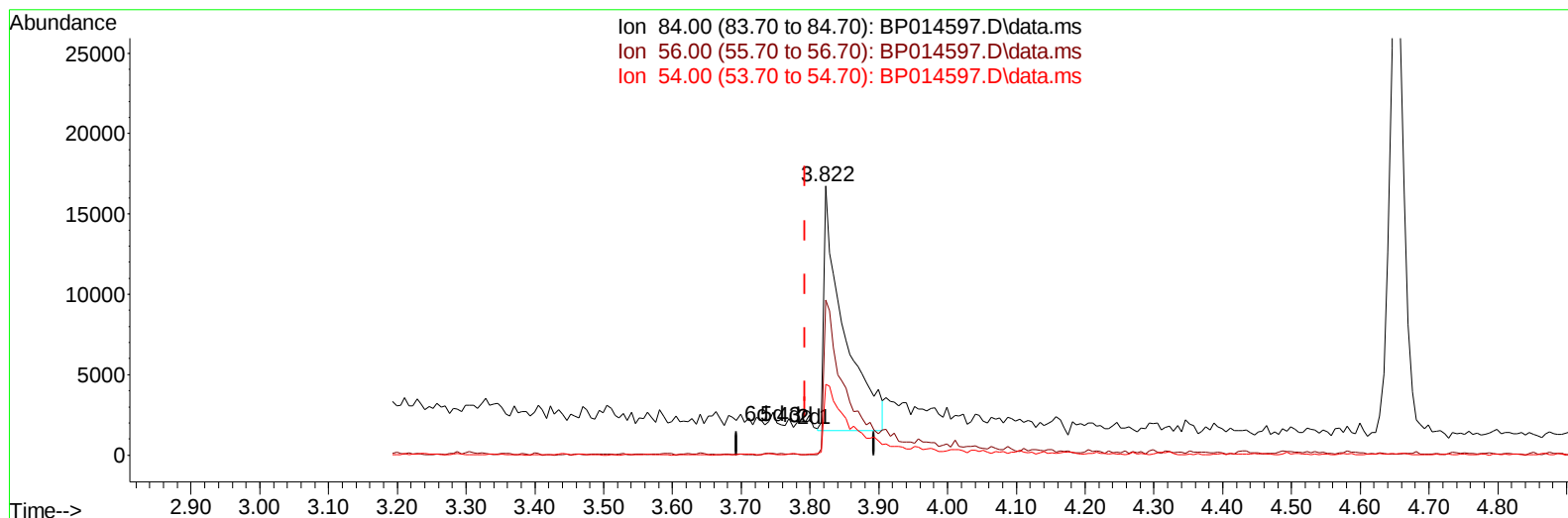
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(4) Pyridine-d5 (S)

3.822min (+ 0.029) 3.20 ng/u1 m

response 30131

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	57.76
54.00	32.60	26.40
0.00	0.00	0.00

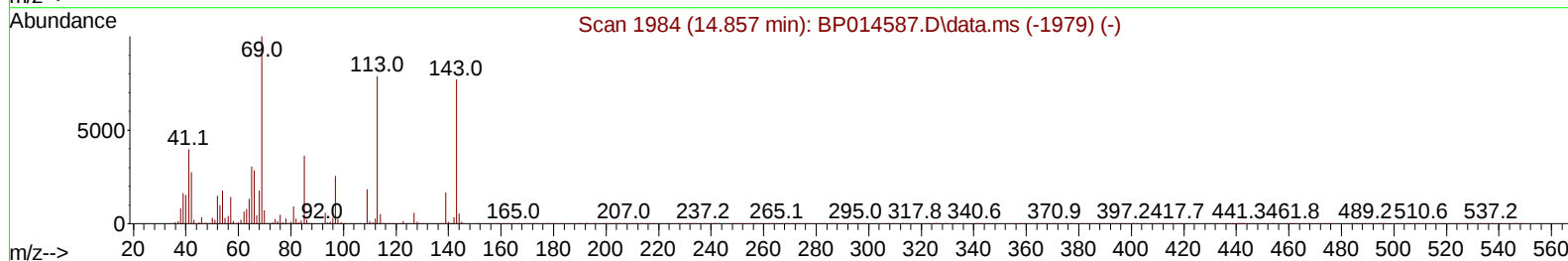
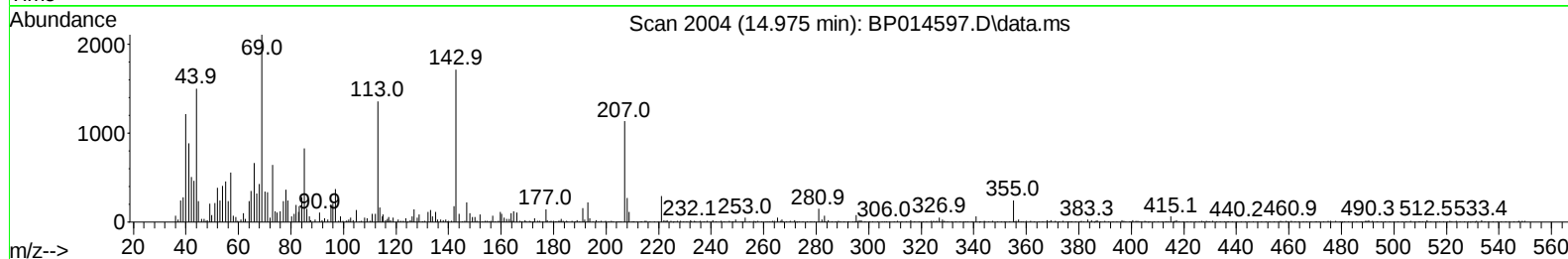
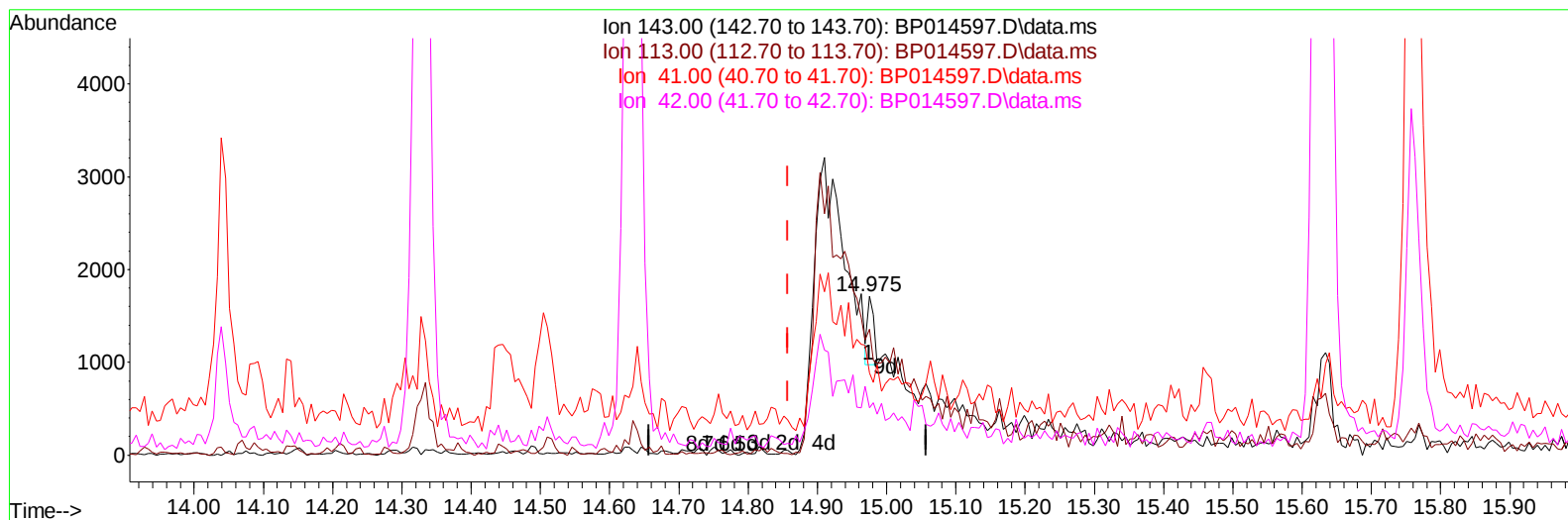
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Instrument :
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TIC: BP014597.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.118) 0.09 ng/u1

response 449

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	79.28
41.00	44.20	51.66
42.00	31.50	29.77

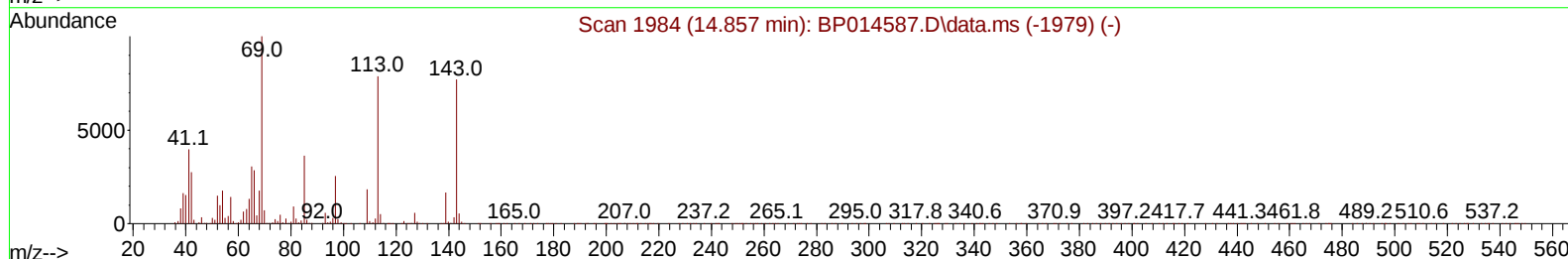
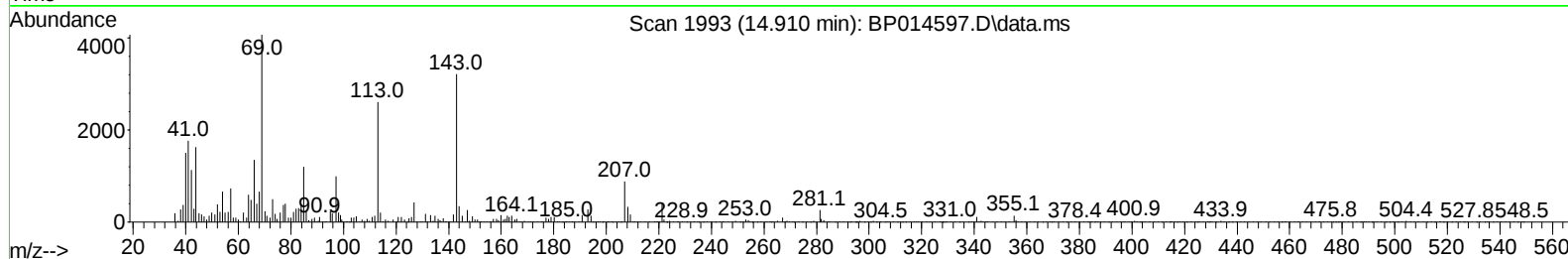
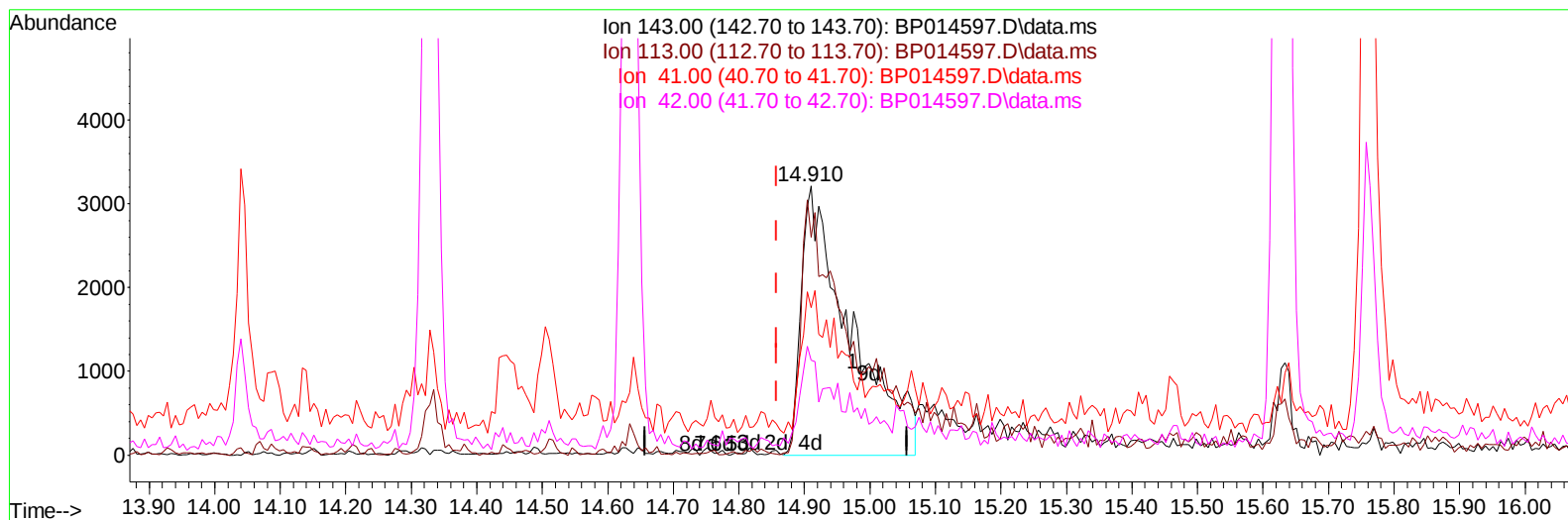
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ALS Vial : 48 Sample Multiplier: 1

Instrument :
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TIC: BP014597.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.053) 3.20 ng/u1 m

response 16753

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	81.03
41.00	44.20	54.91#
42.00	31.50	35.22

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	142595	20.000	ng/ul	0.00
20) Naphthalene-d8	10.804	136	633148	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	427002	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	922213	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	664560	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	683603	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	14184	3.857	ng/uL	0.00
4) Pyridine-d5	3.822	84	30131m	3.198	ng/ul	0.03
7) Phenol-d5	7.158	99	62287	4.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.311	67	224661	26.626	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	189632	19.951	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	126810	11.954	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	125122	24.479	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	137817	23.547	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	223648	20.982	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	164684	11.729	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	965184	27.878	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	1012027	26.343	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	16753m	3.203	ng/ul	0.05
60) Fluorene-d10	15.628	176	812501	28.141	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	151039	22.258	ng/ul	0.00
73) Anthracene-d10	17.492	188	1301623	29.164	ng/ul	0.00
81) Pyrene-d10	19.728	212	1421339	31.841	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	1086625	29.919	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.405	88	10706	2.649	ng/uL	91

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

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