

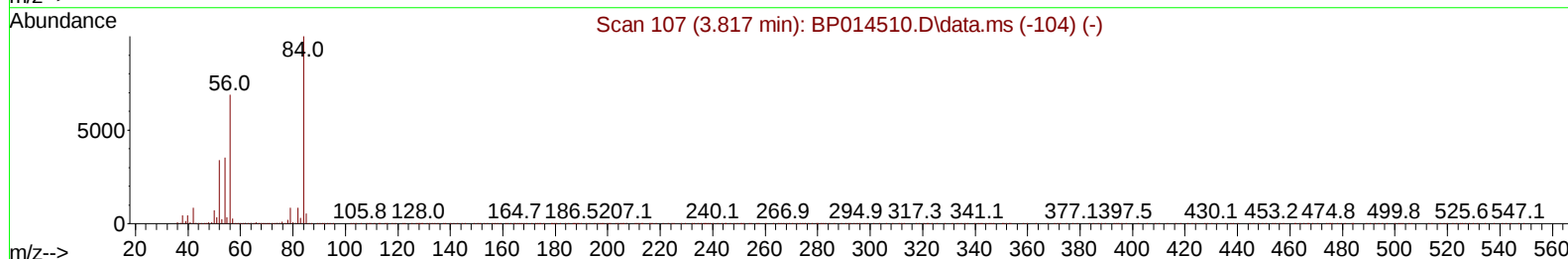
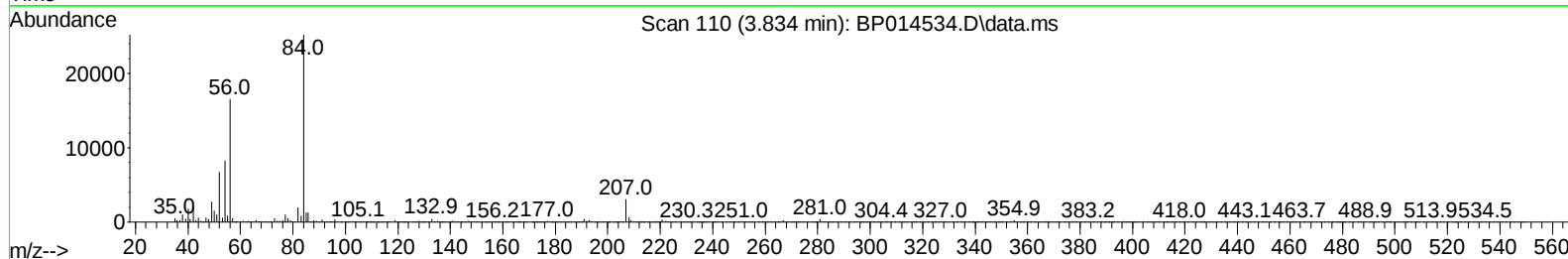
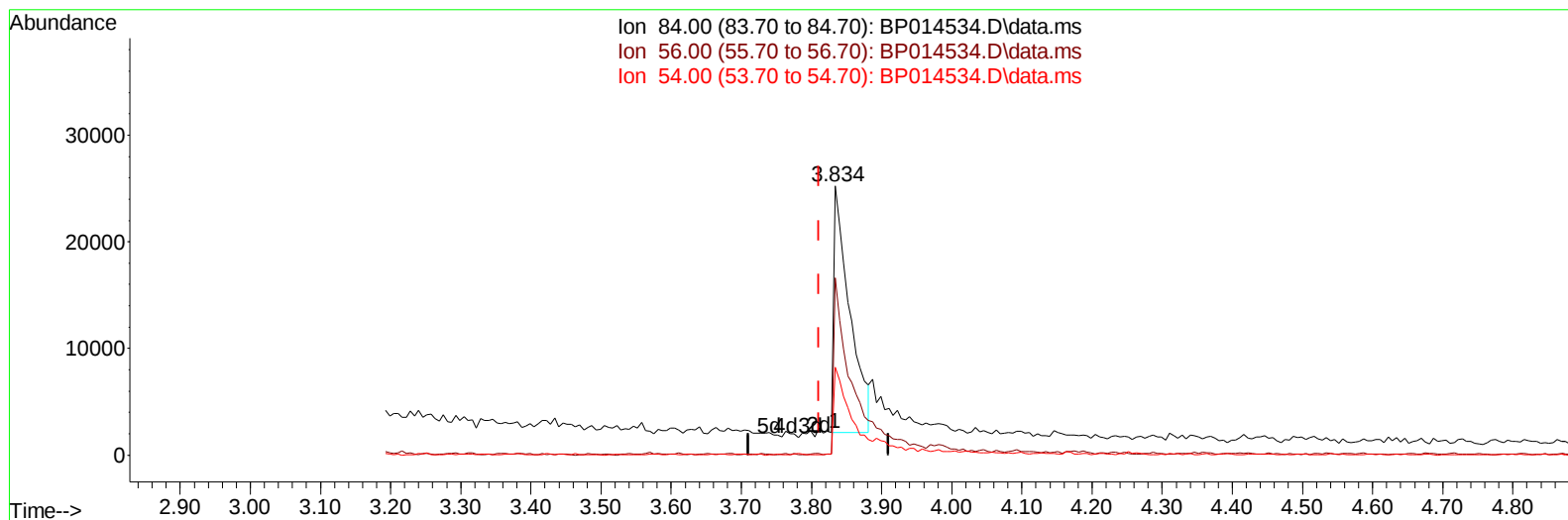
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014534.D
 Acq On : 04 Apr 2023 18:36
 Operator : CG/JU
 Sample : 02158-21
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB(20230331)

Manual IntegrationsAPPROVED

Quant Time: Apr 04 23:46:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 04 23:43:56 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/05/2023
 Supervised By :Jagrut Upadhyay 04/05/2023



TIC: BP014534.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.023) 4.52 ng/u1

response 36573

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.91
54.00	32.60	32.73
0.00	0.00	0.00

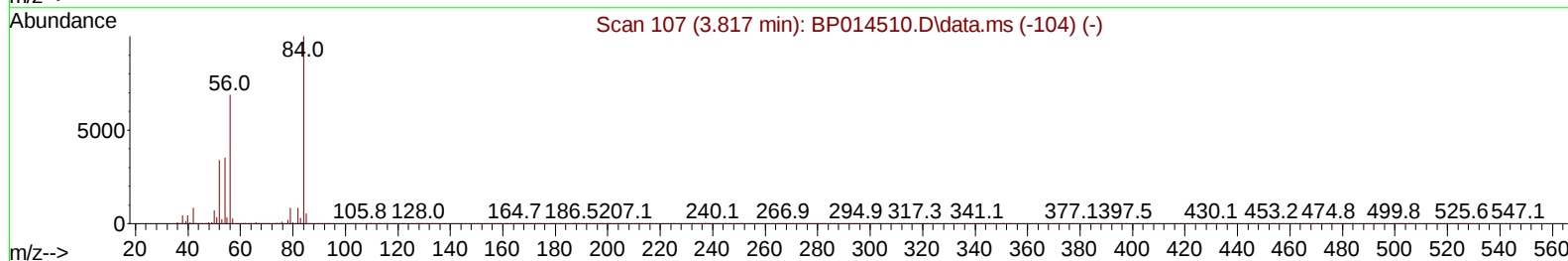
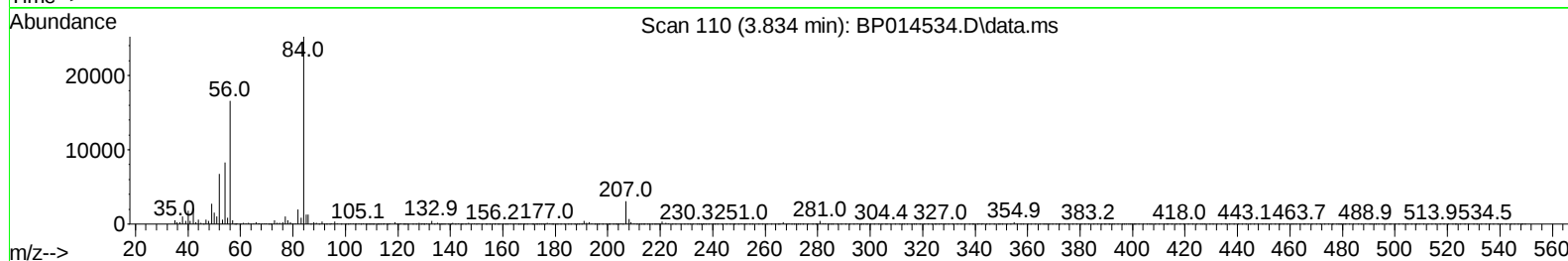
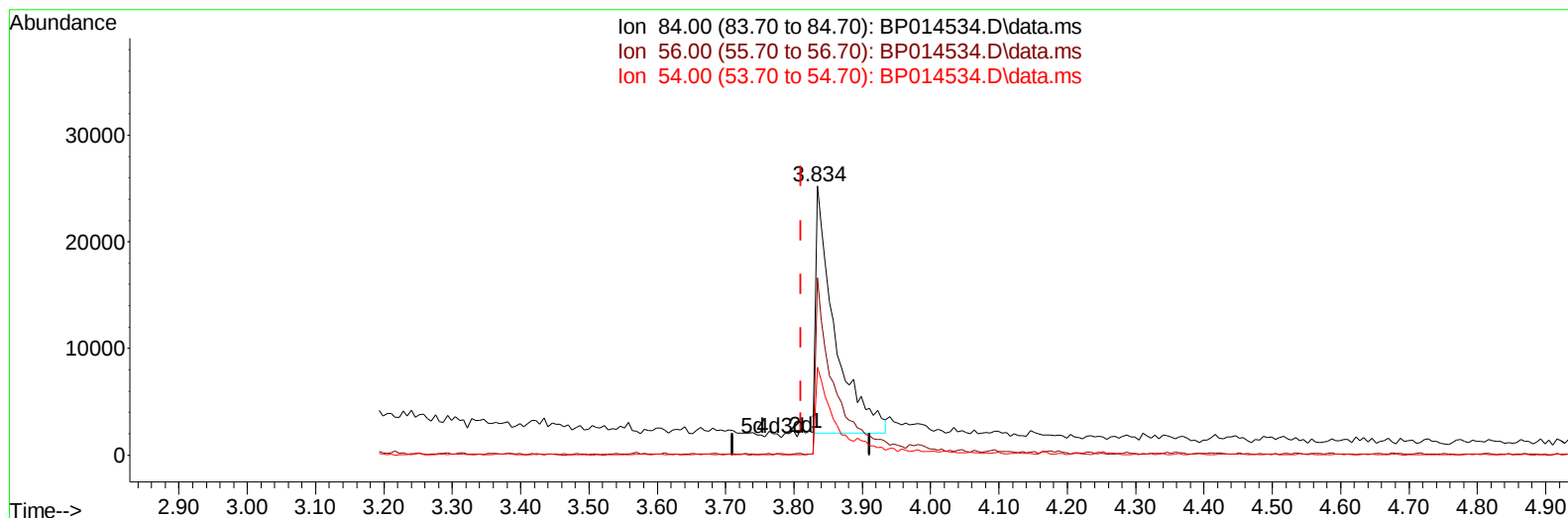
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TIC: BP014534.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.023) 5.52 ng/u1 m

response 44705

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.91
54.00	32.60	32.73
0.00	0.00	0.00

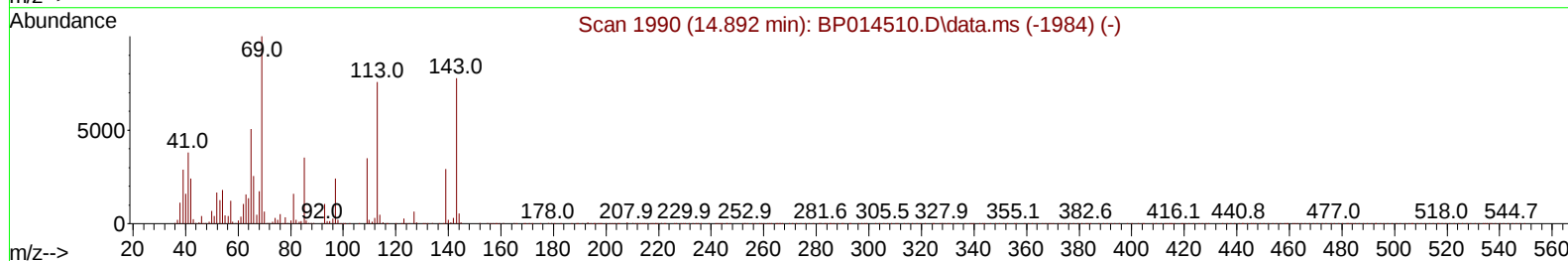
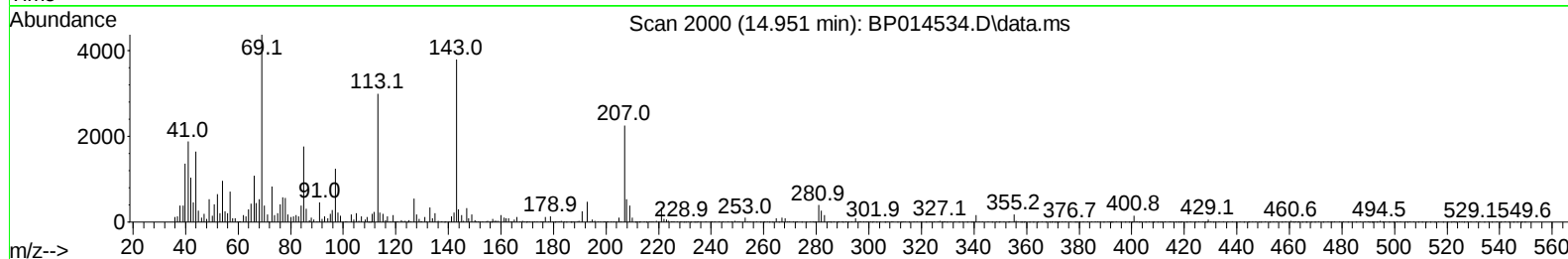
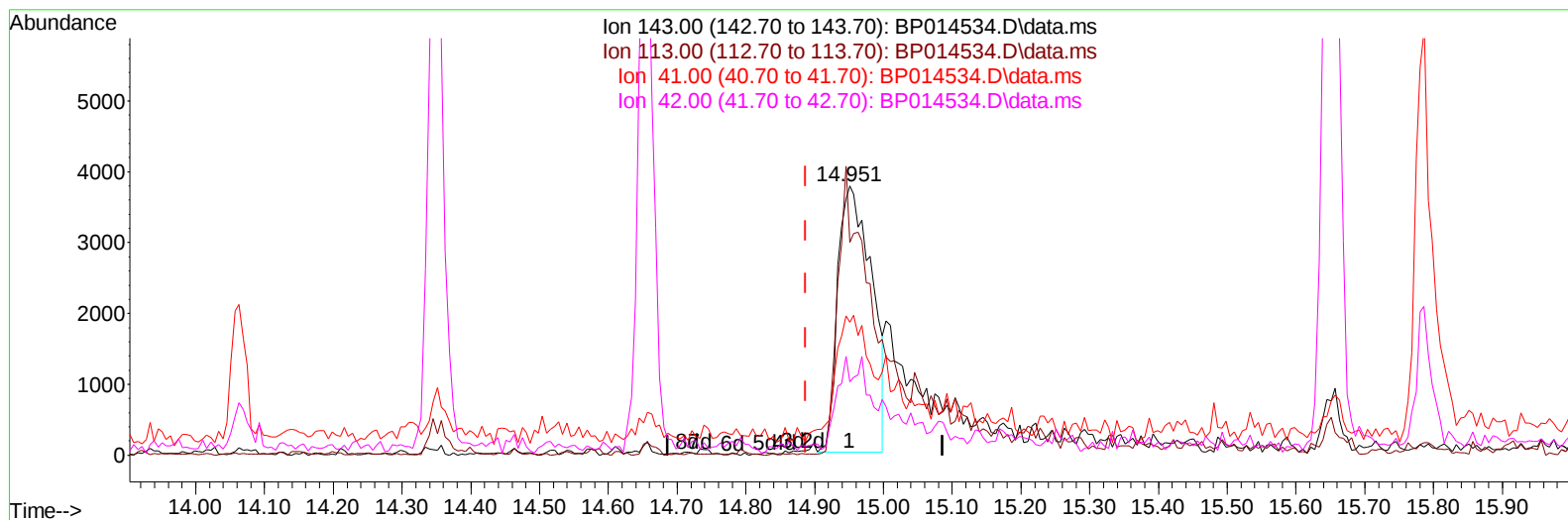
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TIC: BP014534.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.064) 3.04 ng/u1

response 12799

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	78.99
41.00	44.20	49.37
42.00	31.50	27.30

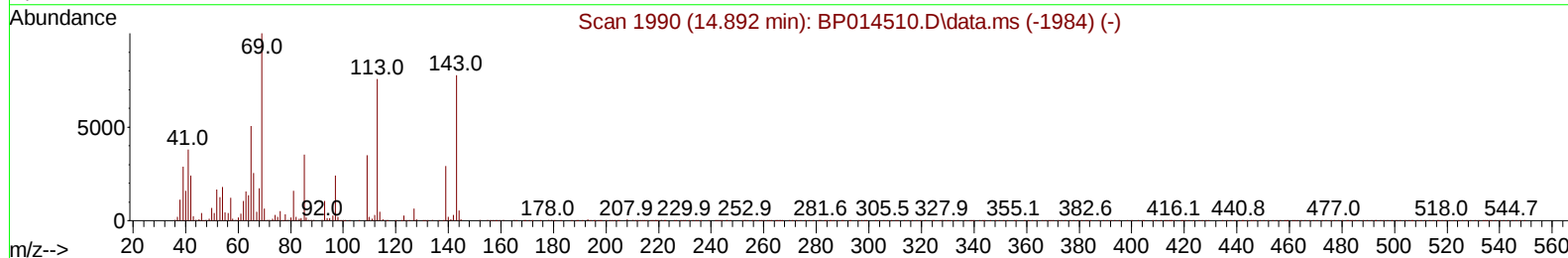
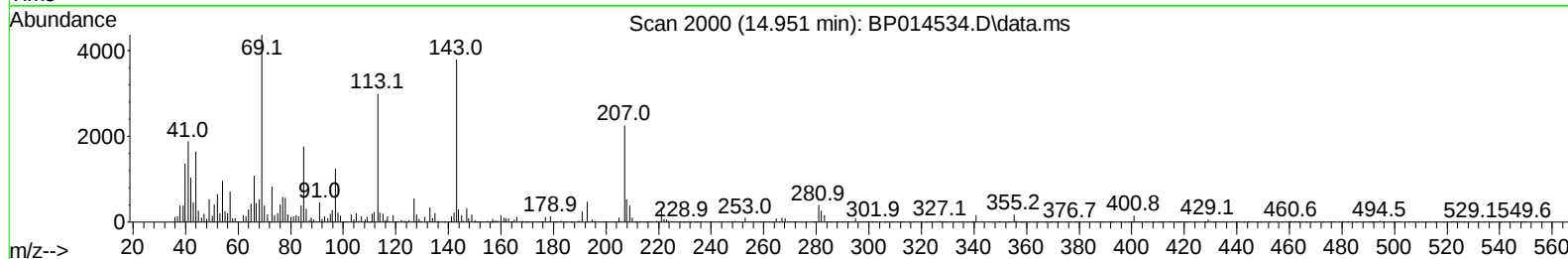
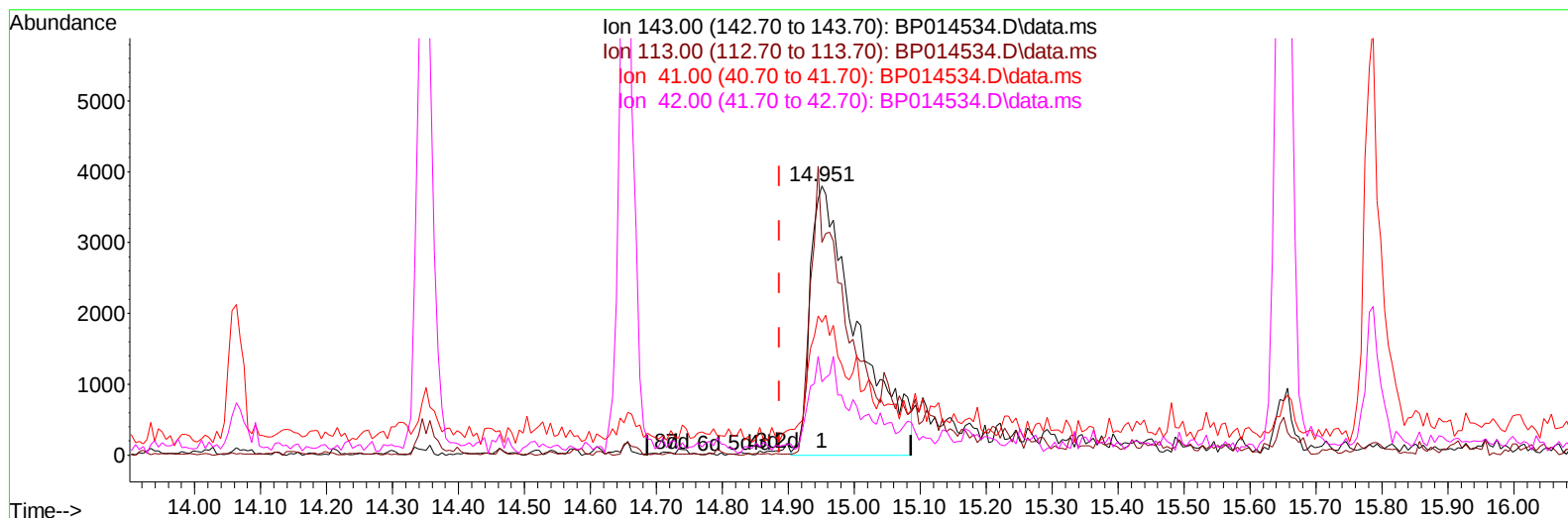
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Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
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TIC: BP014534.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.064) 4.45 ng/ul m

response 18744

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	78.99
41.00	44.20	49.37
42.00	31.50	27.30

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	122547	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	519097	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	344063	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	822645	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	861082	20.000	ng/ul	0.00
88) Perylene-d12	24.015	264	981245	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	15291	4.839	ng/uL	0.00
4) Pyridine-d5	3.834	84	44705m	5.520	ng/ul	0.02
7) Phenol-d5	7.181	99	59733	5.285	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.328	67	192107	26.493	ng/ul	0.00
11) 2-Chlorophenol-d4	7.534	132	169541	20.755	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	114721	12.584	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	109562	26.145	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	116478	24.274	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.457	165	187207	21.422	ng/ul	0.00
31) 4-Chloroaniline-d4	10.975	131	234447	20.367	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	779573	27.944	ng/ul	0.00
49) Acenaphthylene-d8	14.351	160	828276	26.758	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	18744m	4.447	ng/ul	0.06
60) Fluorene-d10	15.651	176	680351	29.244	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.786	200	103055	17.025	ng/ul	0.00
73) Anthracene-d10	17.516	188	1252625	31.463	ng/ul	0.00
81) Pyrene-d10	19.745	212	1567565	27.102	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.851	264	1650258	31.655	ng/ul	0.00

Target Compounds Qvalue

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

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