

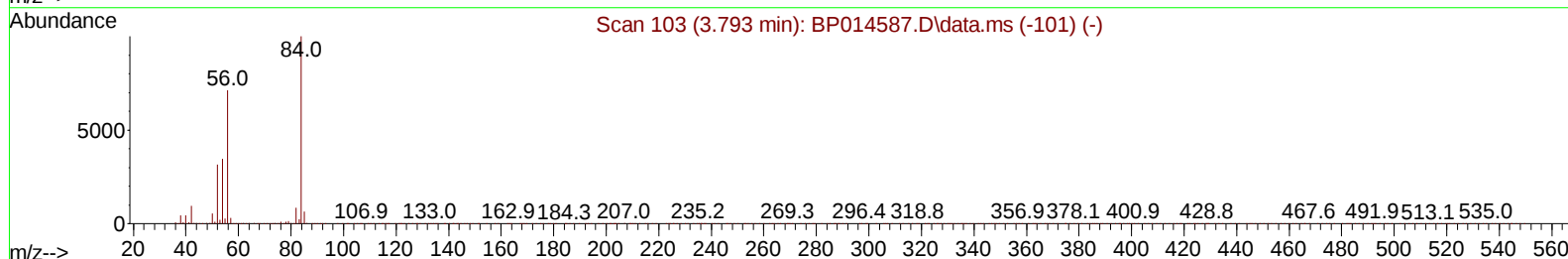
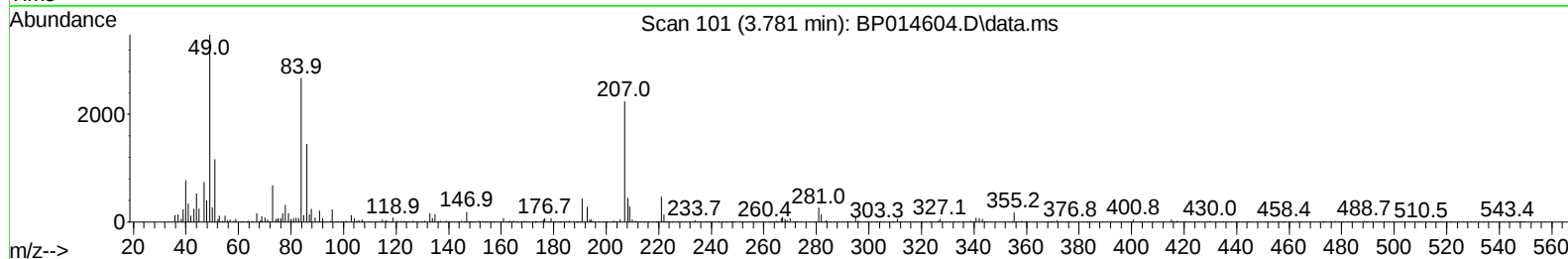
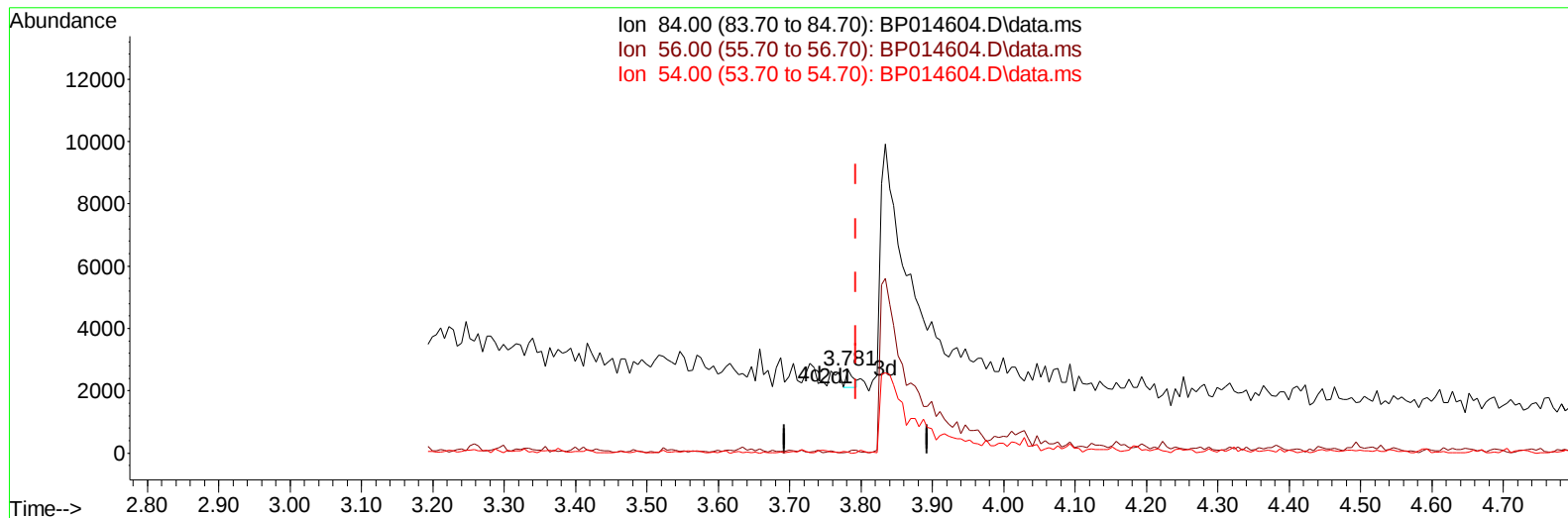
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014604.D
Acq On : 07 Apr 2023 00:37
Operator : CG/JU
Sample : 02225-06
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AD2

Manual IntegrationsAPPROVED

Quant Time: Apr 07 01:23:33 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
Response via : Initial Calibration

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



TIC: BP014604.D\data.ms

(4) Pyridine-d5 (S)

3.781min (-0.012) 0.04 ng/u1

response 409

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	1.23#
54.00	32.60	0.97#
0.00	0.00	0.00

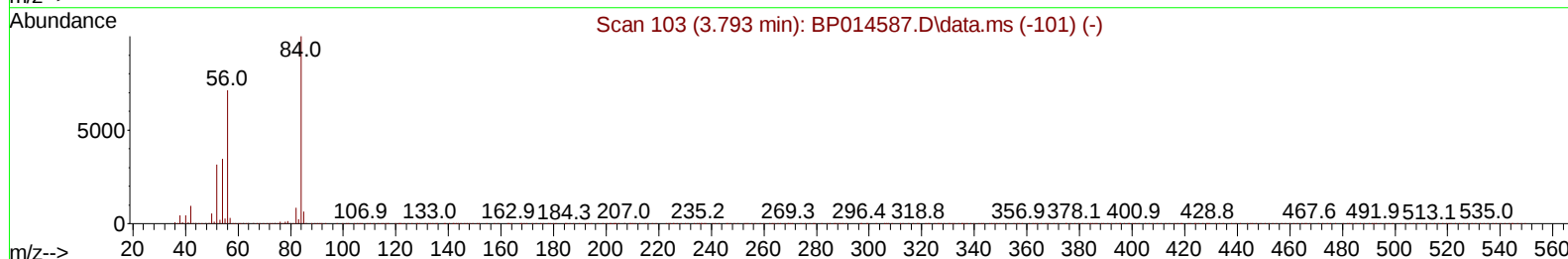
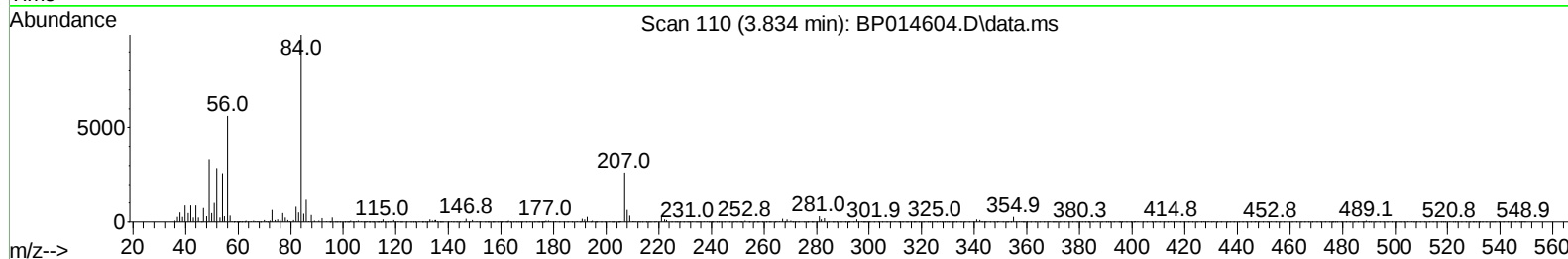
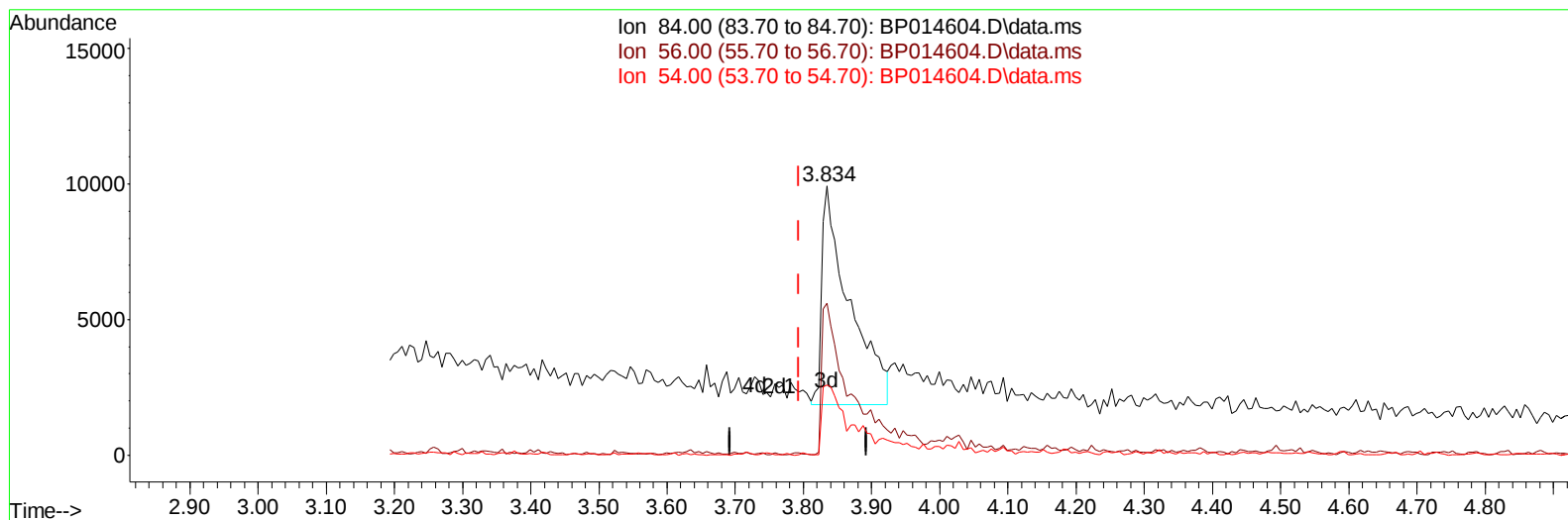
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(4) Pyridine-d5 (S)

3.834min (+ 0.041) 2.17 ng/ul m

response 22665

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	56.50
54.00	32.60	25.97#
0.00	0.00	0.00

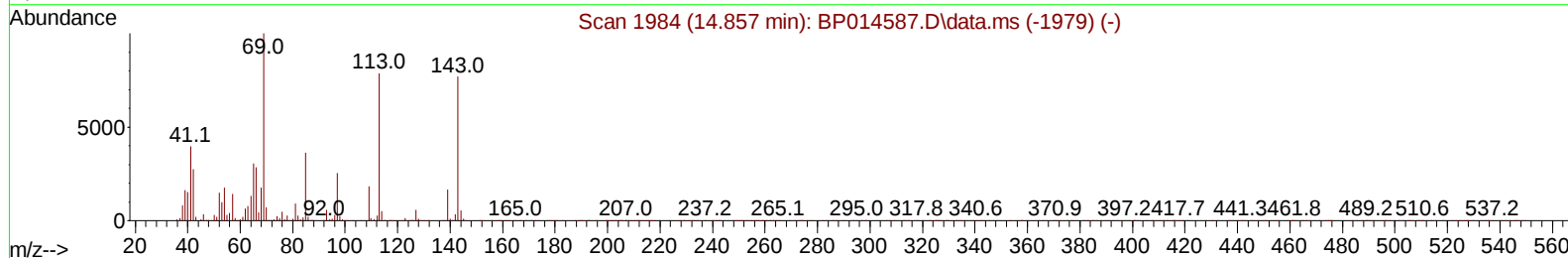
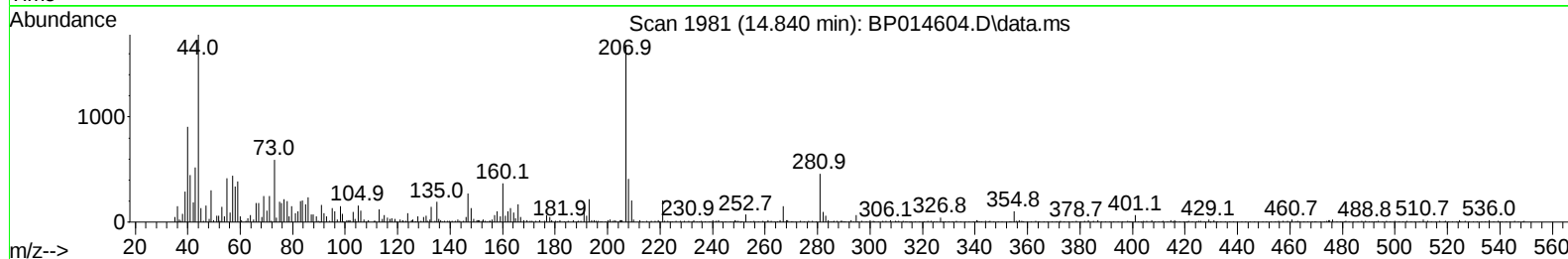
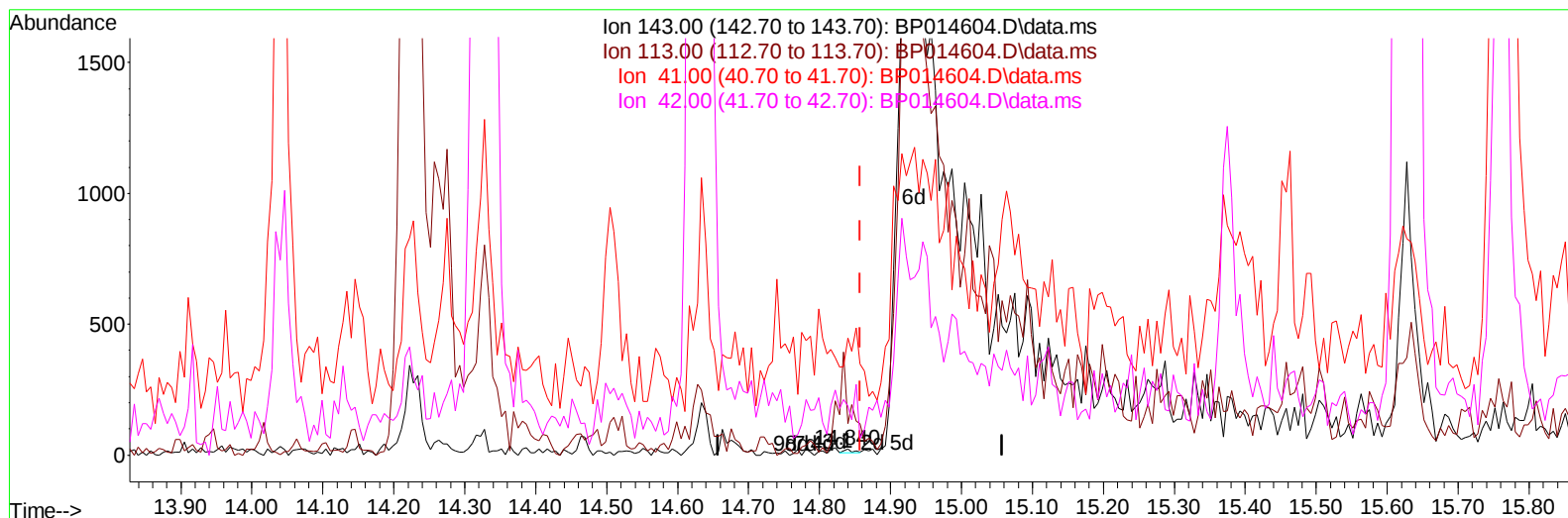
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ALS Vial : 55 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014604.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.840min (-0.018) 0.00 ng/u1

response 11

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	324.32#
41.00	44.20	1205.41#
42.00	31.50	508.11#

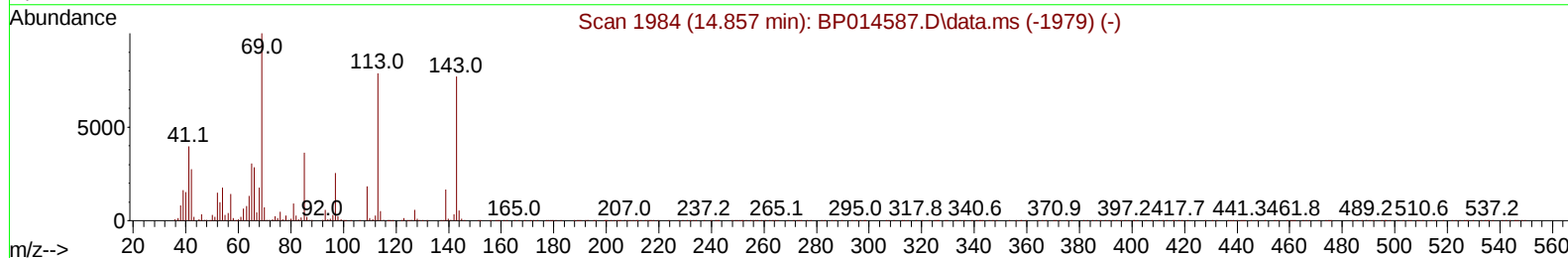
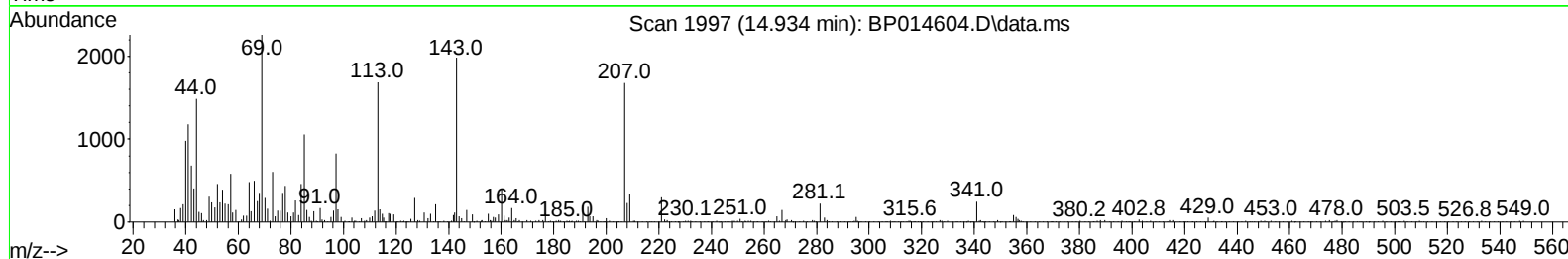
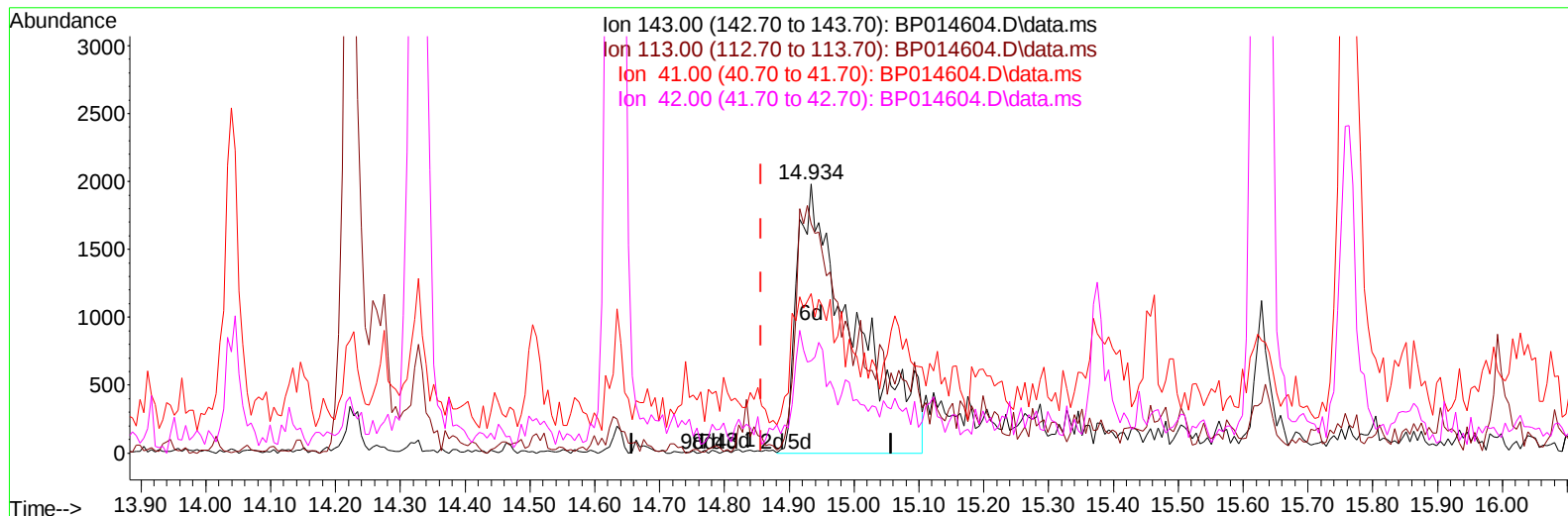
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Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
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TIC: BP014604.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.934min (+ 0.077) 2.01 ng/ul m

response 12004

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	85.20
41.00	44.20	59.39#
42.00	31.50	34.29

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 ALS Vial : 55 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	158039	20.000	ng/ul	0.00
20) Naphthalene-d8	10.799	136	722662	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.634	164	487746	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	1078486	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	809202	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	804187	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	13874	3.404	ng/uL	0.00
4) Pyridine-d5	3.834	84	22665m	2.170	ng/ul	0.04
7) Phenol-d5	7.164	99	40188	2.757	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.311	67	184924	19.775	ng/ul	0.00
11) 2-Chlorophenol-d4	7.511	132	131421	12.475	ng/ul	0.00
15) 4-Methylphenol-d8	8.705	113	87909	7.477	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128	98895	16.952	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	94539	14.152	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	160090	13.159	ng/ul	0.00
31) 4-Chloroaniline-d4	10.958	131	196506	12.262	ng/ul	0.00
46) Dimethylphthalate-d6	14.040	166	835342	21.123	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	842466	19.198	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143	12004m	2.009	ng/ul	0.08
60) Fluorene-d10	15.628	176	688606	20.880	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	105543	13.300	ng/ul	0.00
73) Anthracene-d10	17.492	188	1165337	22.327	ng/ul	0.00
81) Pyrene-d10	19.722	212	1331694	24.500	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	980518	22.949	ng/ul	0.00

Target Compounds Qvalue

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

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