

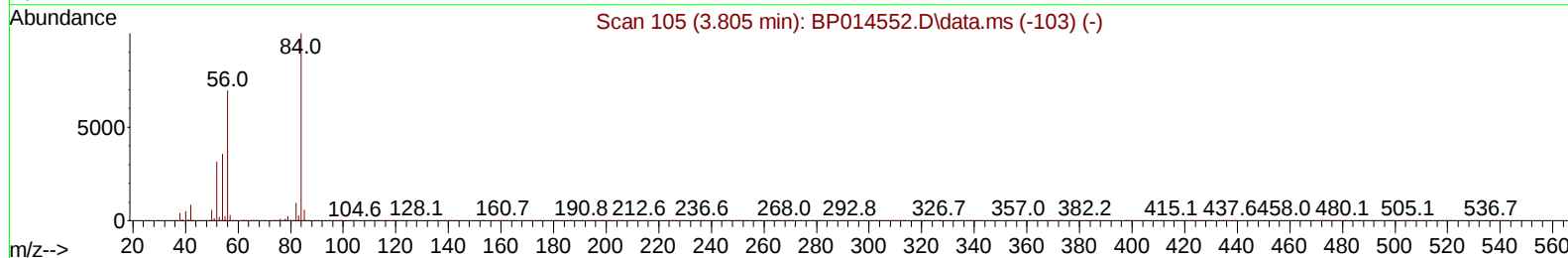
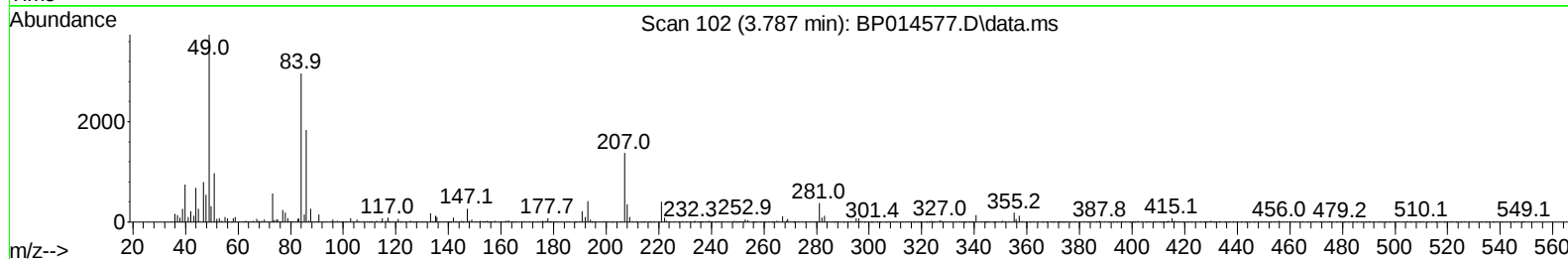
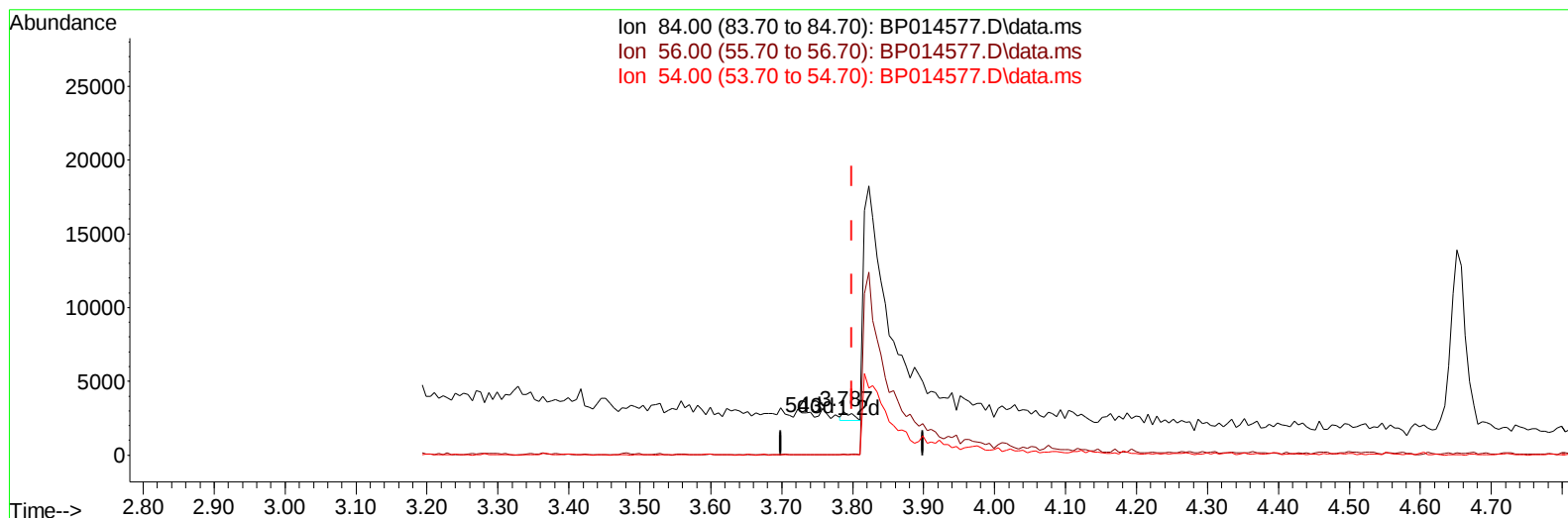
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014577.D
 Acq On : 06 Apr 2023 03:15
 Operator : CG/JU
 Sample : 02194-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHDW0

Manual IntegrationsAPPROVED

Quant Time: Apr 06 03:56:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 06 00:55:13 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/06/2023
 Supervised By :Jagrut Upadhyay 04/06/2023



TIC: BP014577.D\data.ms

(4) Pyridine-d5 (S)

3.787min (-0.012) 0.06 ng/u1

response 548

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	2.56#
54.00	32.60	0.74#
0.00	0.00	0.00

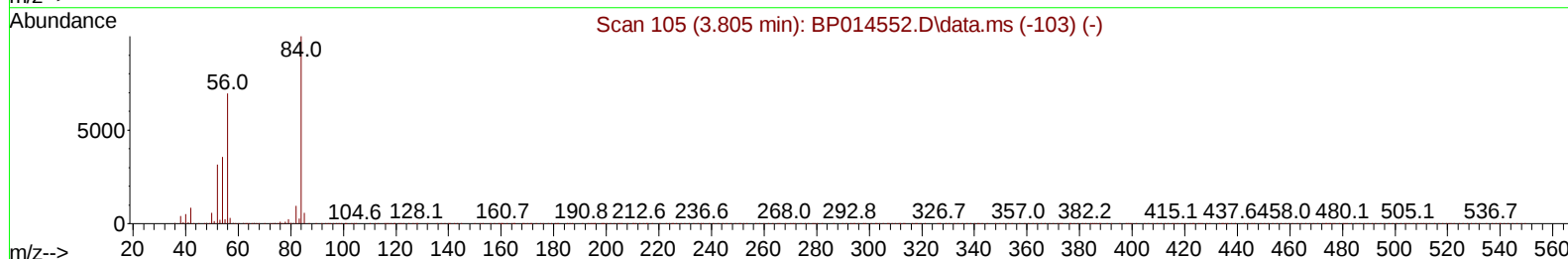
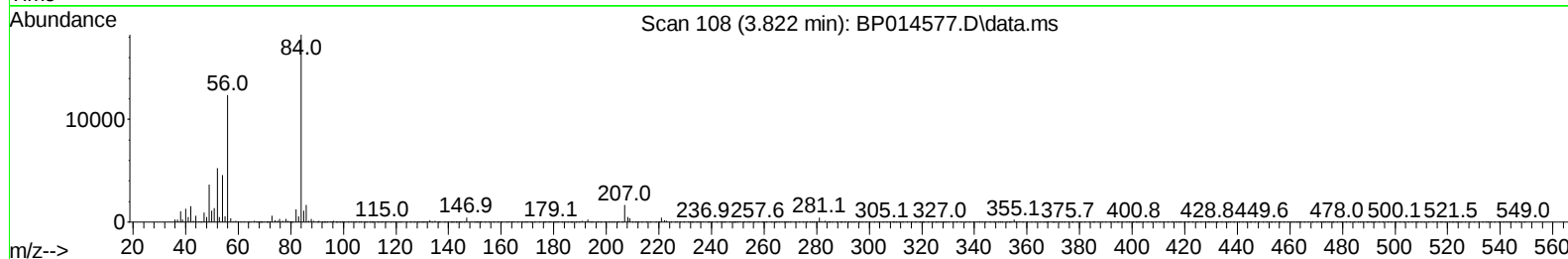
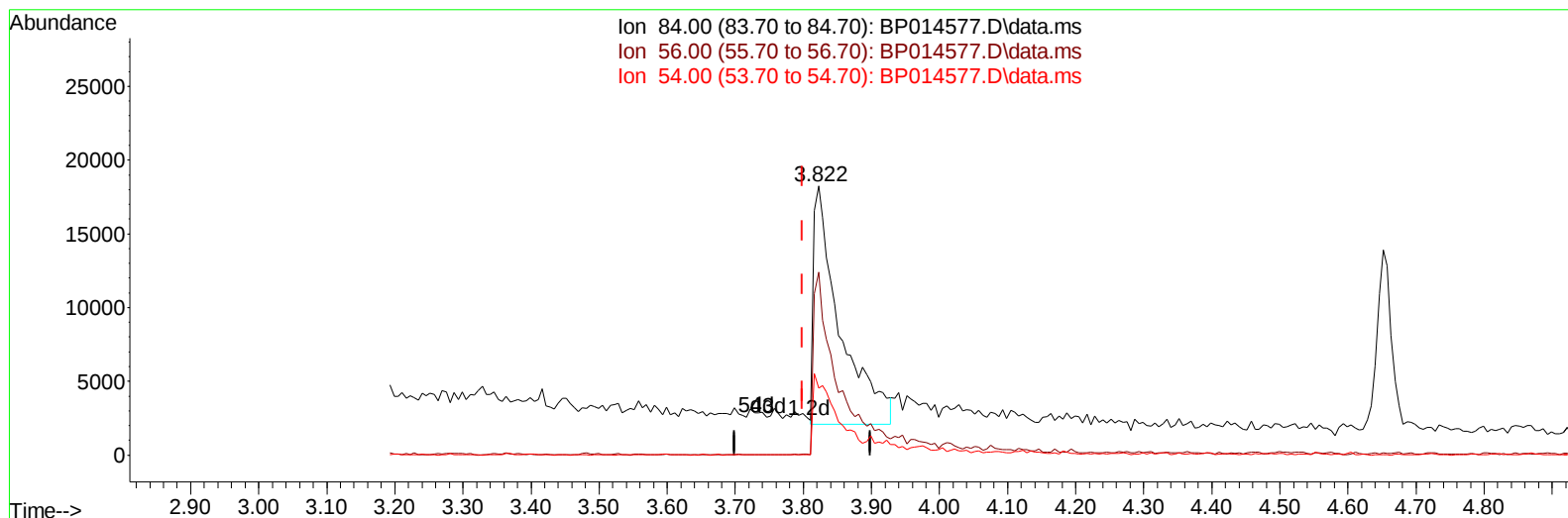
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(4) Pyridine-d5 (S)

3.822min (+ 0.023) 5.10 ng/u1 m

response 43183

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	67.95
54.00	32.60	24.93#
0.00	0.00	0.00

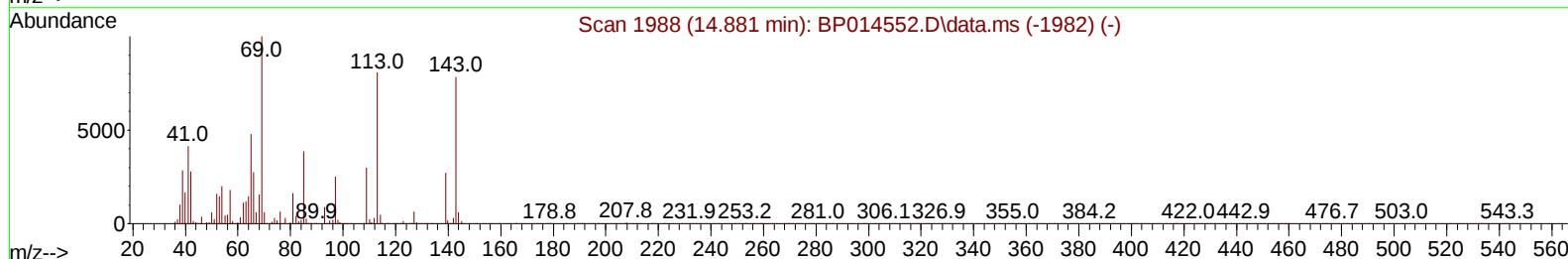
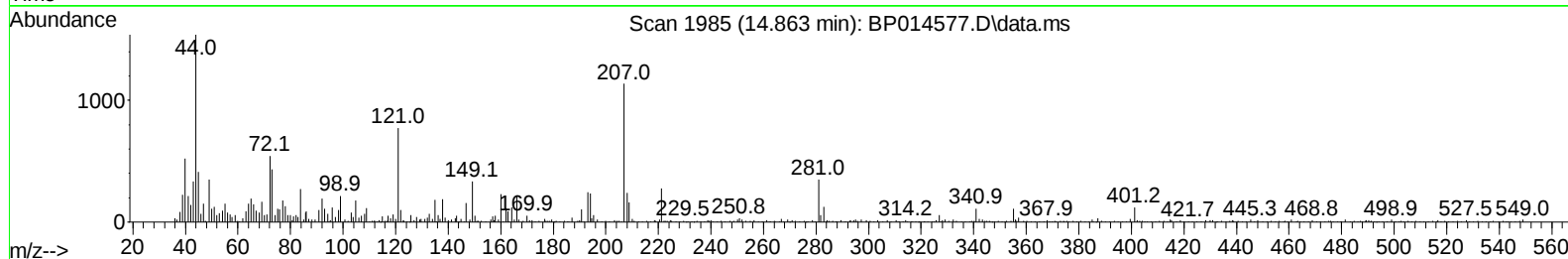
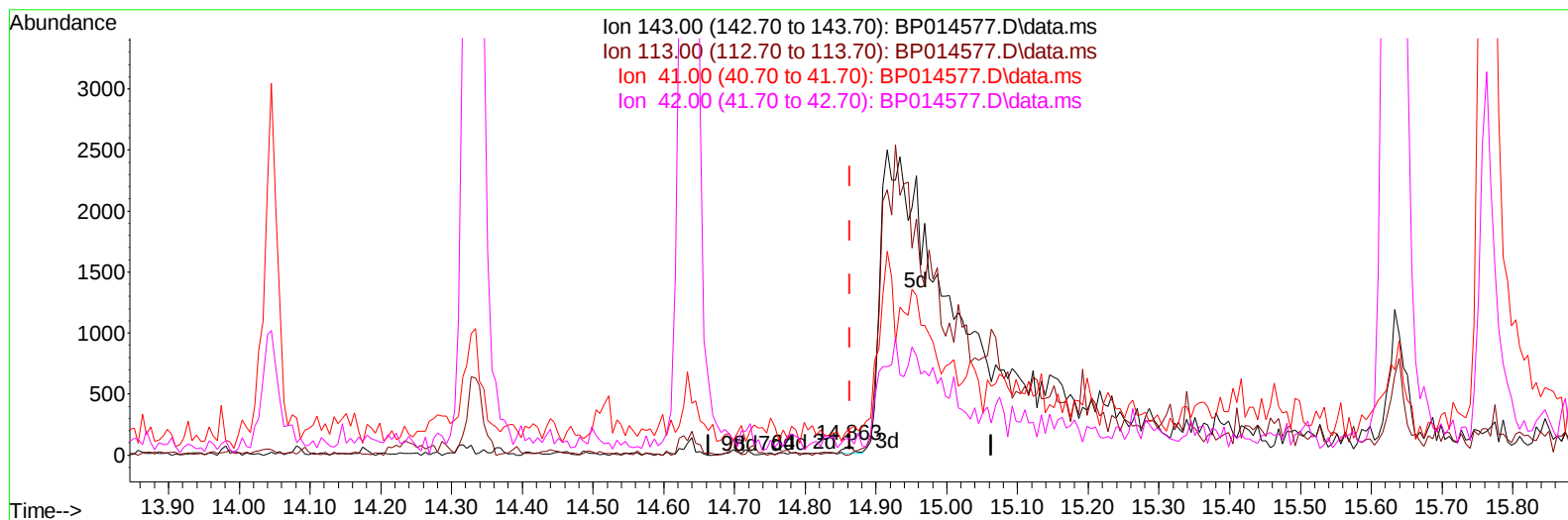
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 Sample : 02194-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP014577.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.863min (-0.000) 0.01 ng/u1

response 62

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	26.19#
41.00	44.20	254.76#
42.00	31.50	167.86#

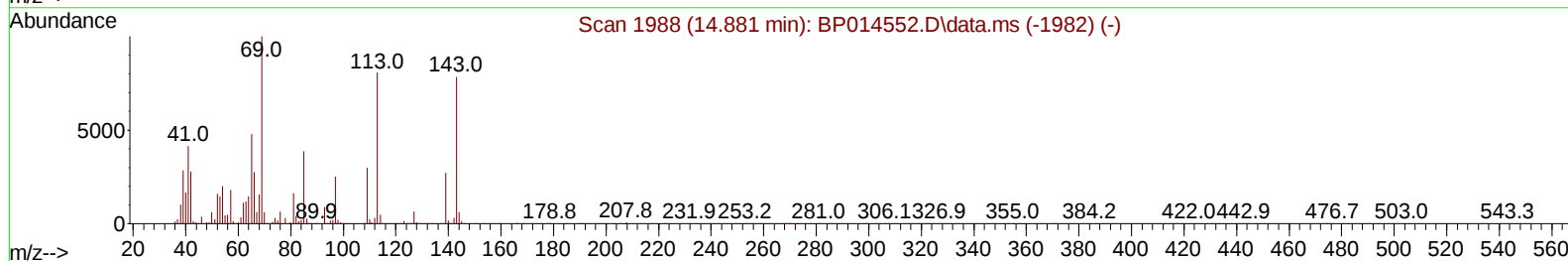
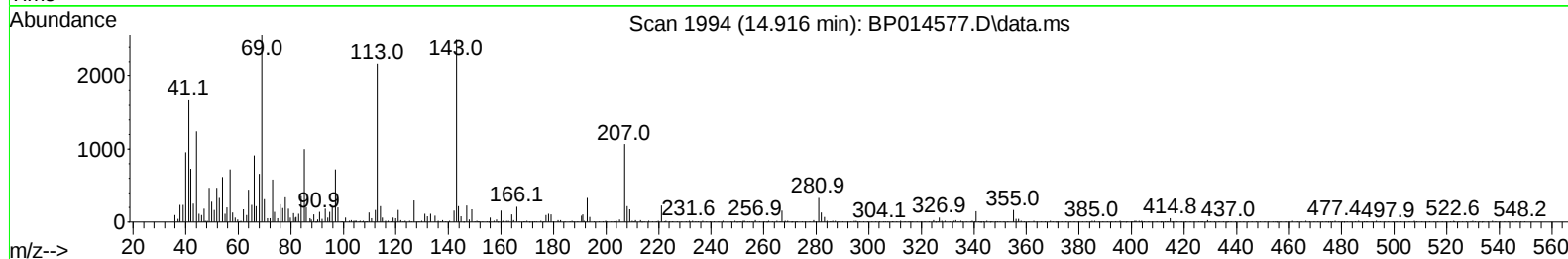
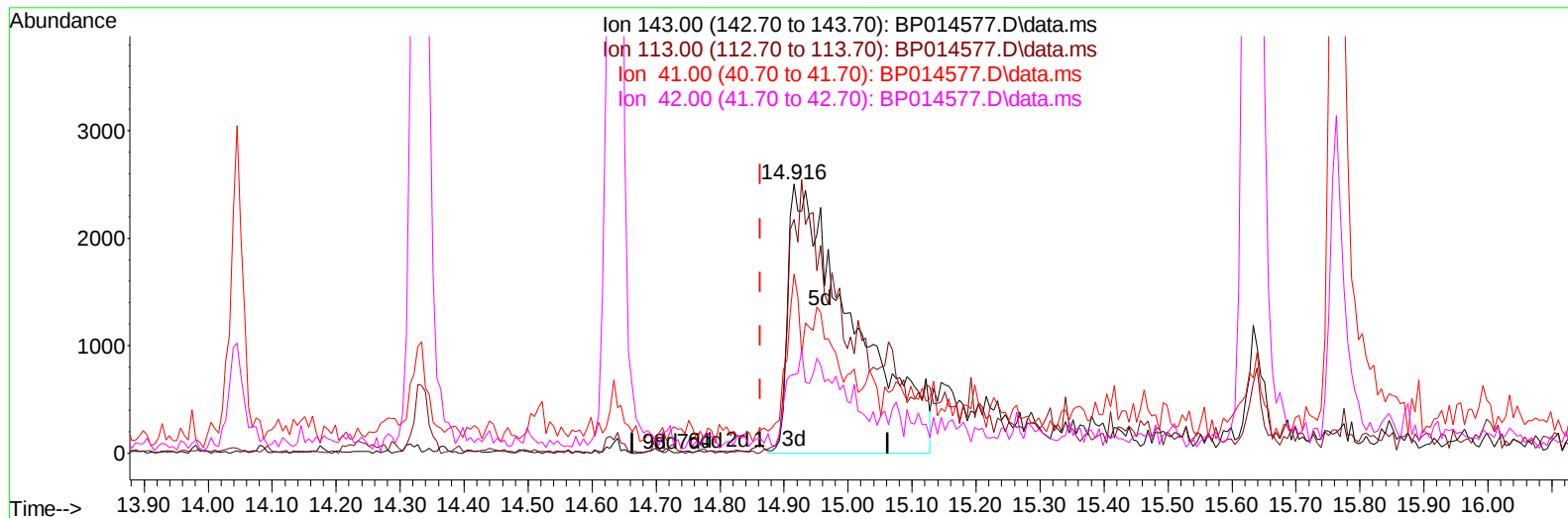
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ALS Vial : 28 Sample Multiplier: 1

Instrument :
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TIC: BP014577.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.916min (+ 0.053) 3.67 ng/ul m

response 17706

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	86.98
41.00	44.20	66.68#
42.00	31.50	29.09

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	128164	20.000	ng/ul	0.00
20) Naphthalene-d8	10.804	136	560255	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	393956	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	899745	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	767995	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	761031	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	12923	3.910	ng/uL	0.00
4) Pyridine-d5	3.822	84	43183m	5.099	ng/ul	0.02
7) Phenol-d5	7.163	99	60074	5.082	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.310	67	228136	30.082	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	186325	21.810	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	122779	12.878	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	125173	27.676	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	132927	25.667	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	206291	21.872	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	194299	15.639	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	1002253	31.377	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	1032675	29.136	ng/ul	0.00
54) 4-Nitrophenol-d4	14.916	143	17706m	3.669	ng/ul	0.05
60) Fluorene-d10	15.634	176	835546	31.367	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	151080	22.820	ng/ul	0.00
73) Anthracene-d10	17.498	188	1444555	33.175	ng/ul	0.00
81) Pyrene-d10	19.727	212	1721714	33.376	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	1362803	33.705	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.405	88	7583	2.088	ng/uL	96

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

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