

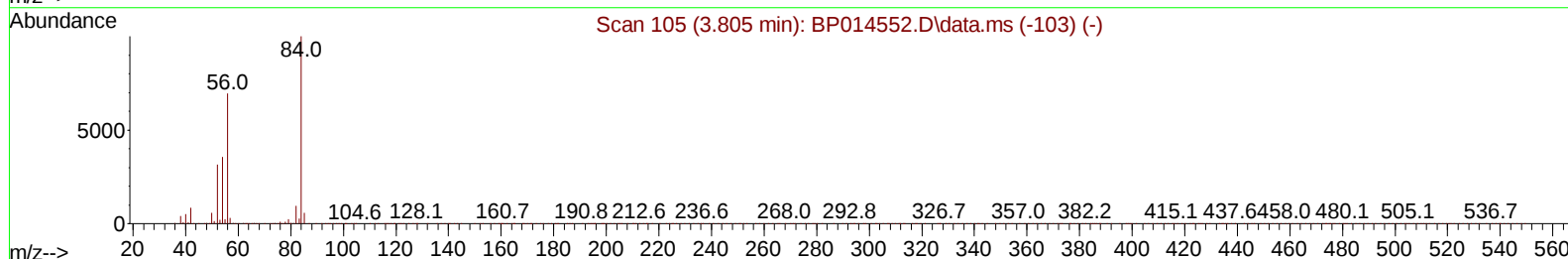
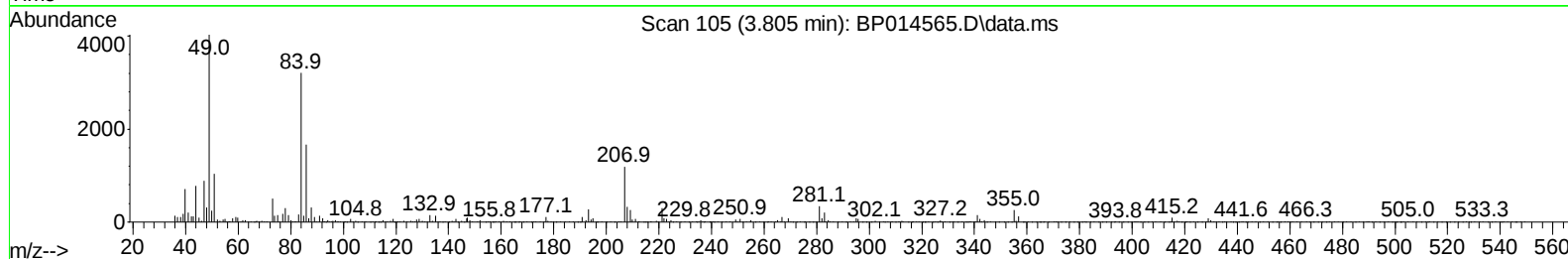
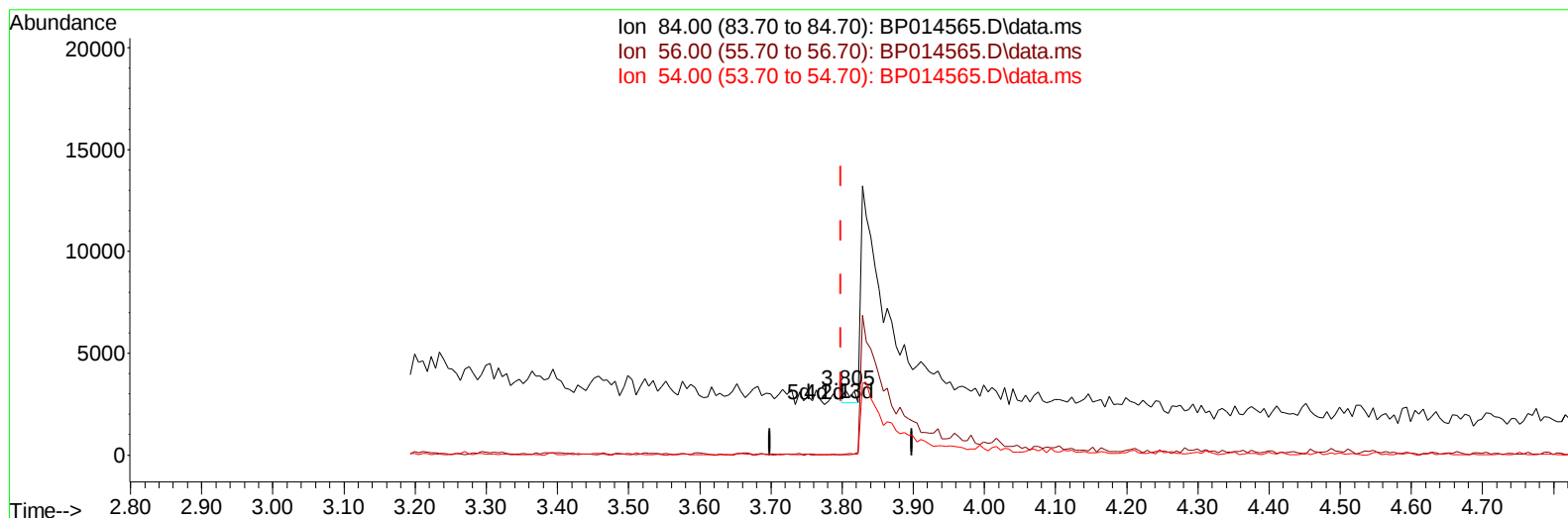
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014565.D
 Acq On : 05 Apr 2023 19:25
 Operator : CG/JU
 Sample : 02122-15
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB9C9

Manual IntegrationsAPPROVED

Quant Time: Apr 06 00:58:59 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: BP014565.D\data.ms

(4) Pyridine-d5 (S)

3.805min (+ 0.006) 0.06 ng/u1

response 530

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.72#
54.00	32.60	1.84#
0.00	0.00	0.00

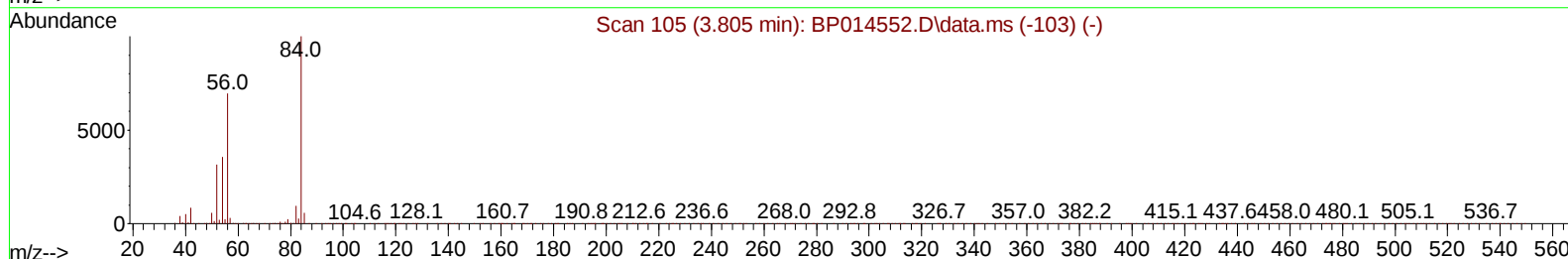
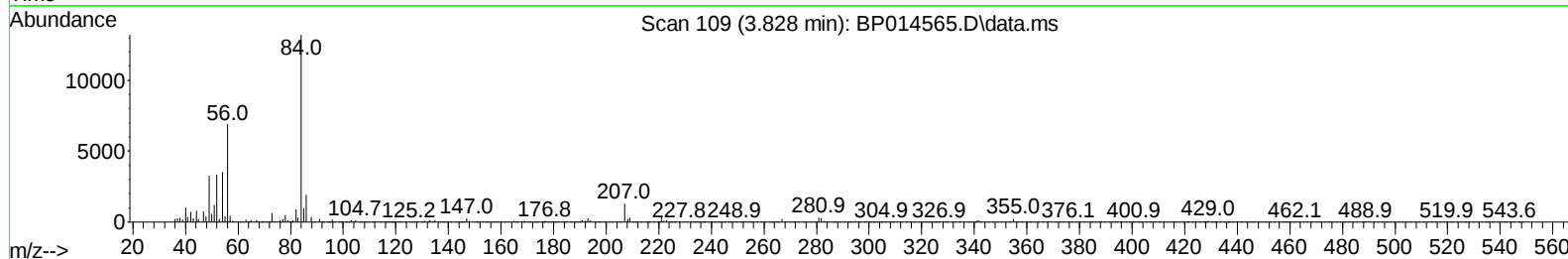
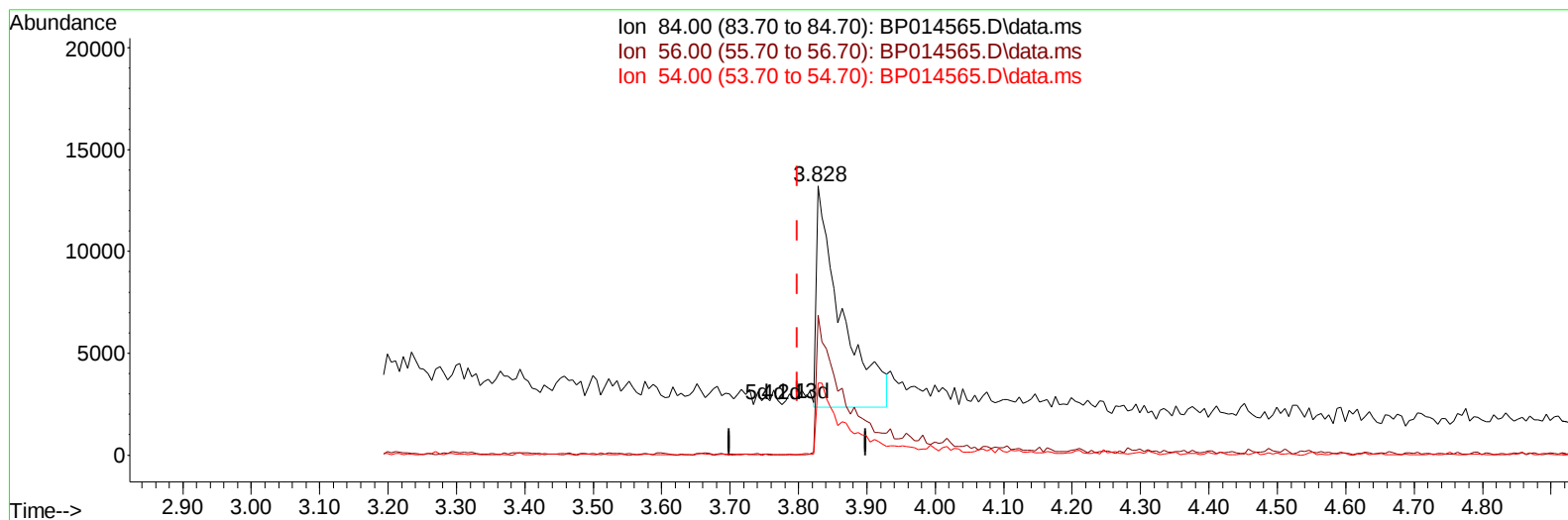
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(4) Pyridine-d5 (S)

3.828min (+ 0.029) 3.10 ng/ul m

response 26996

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	52.10#
54.00	32.60	26.89
0.00	0.00	0.00

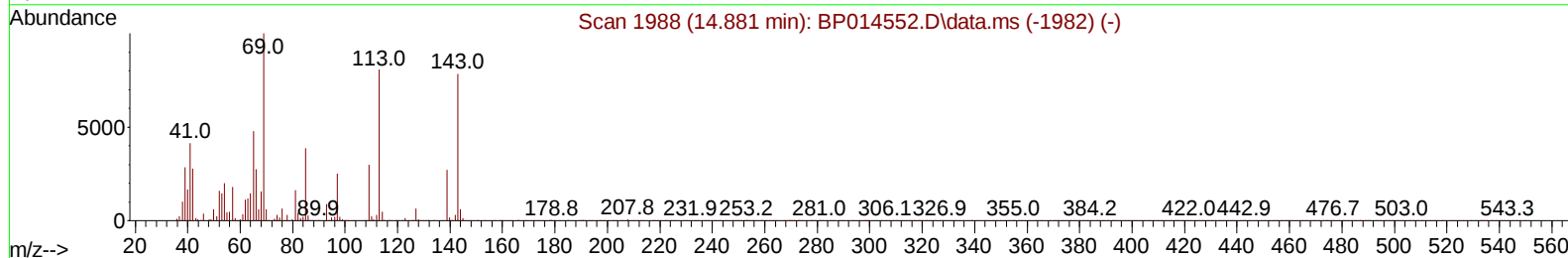
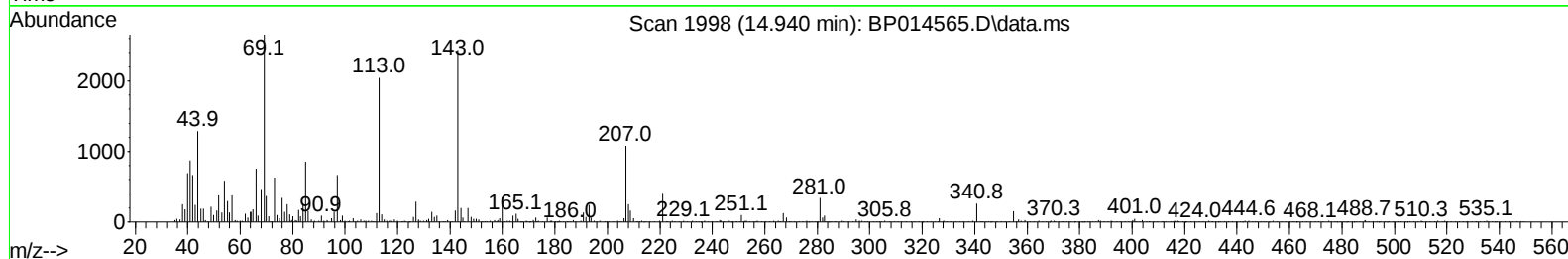
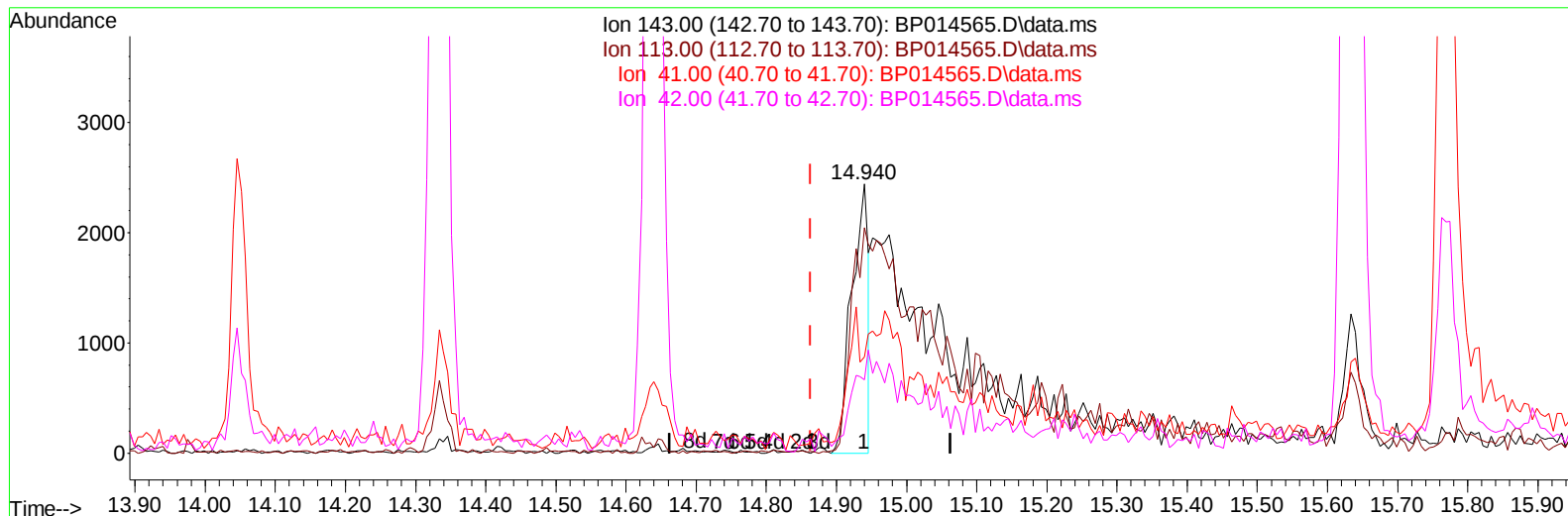
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP014565.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.940min (+ 0.077) 0.83 ng/u1

response 4028

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.78
41.00	44.20	35.71
42.00	31.50	27.35

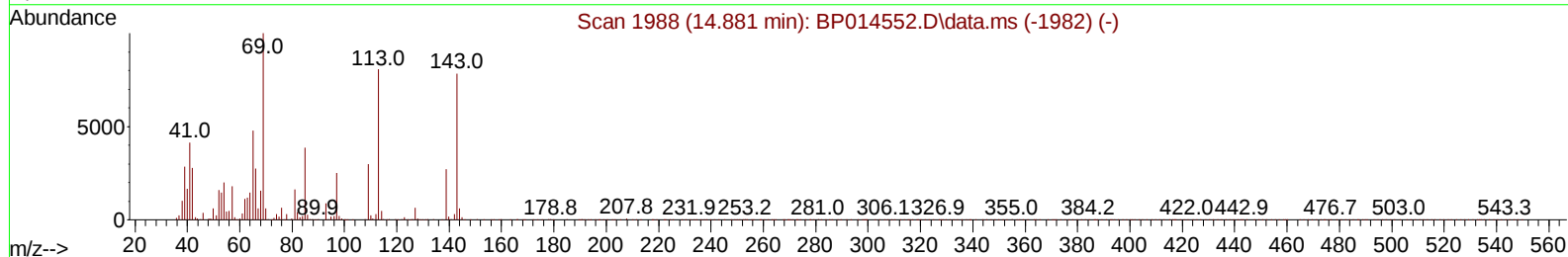
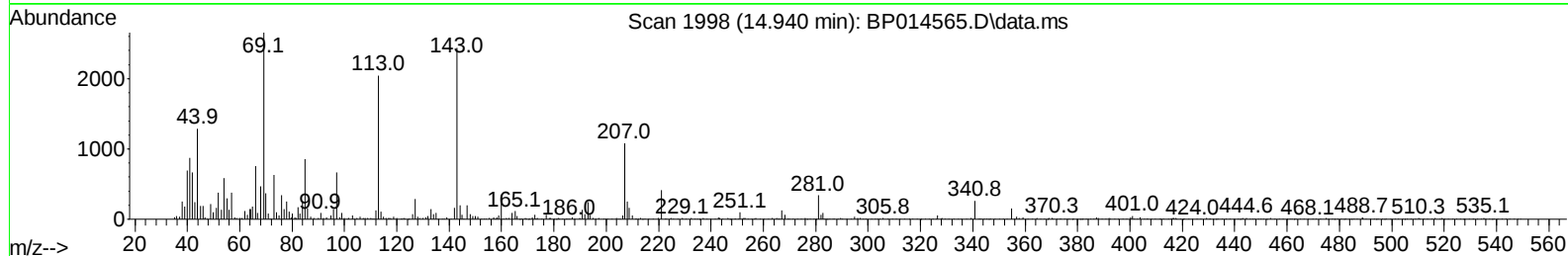
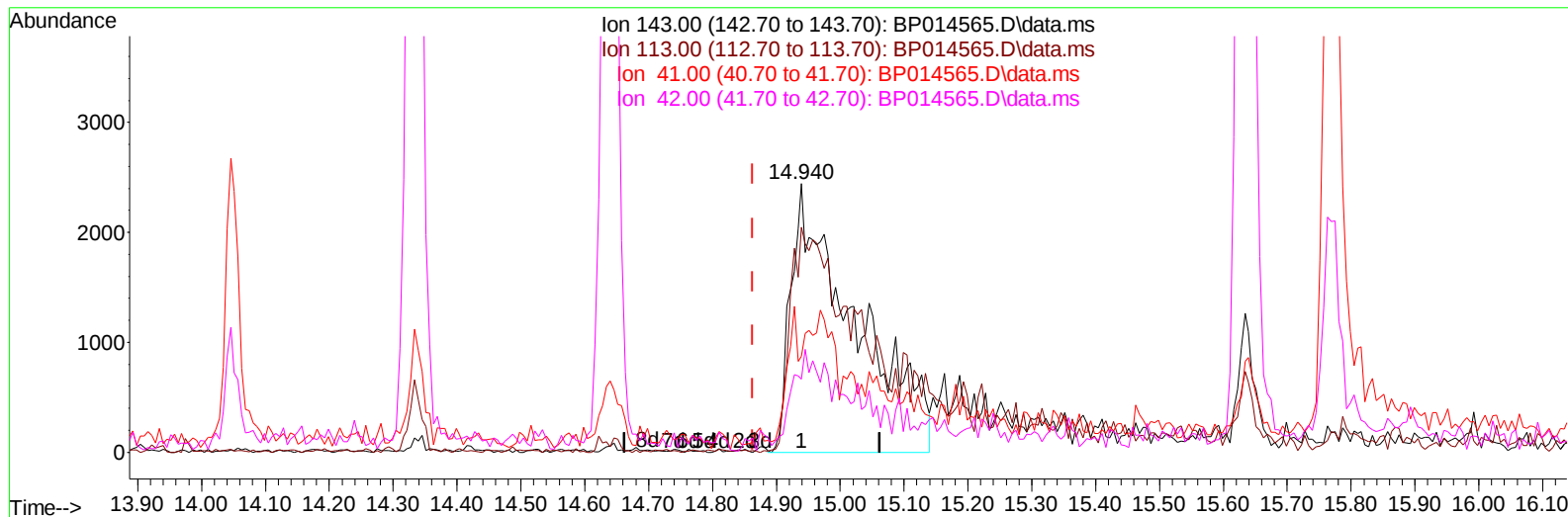
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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TIC: BP014565.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.940min (+ 0.077) 3.49 ng/ul m

response 16936

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	83.78
41.00	44.20	35.71
42.00	31.50	27.35

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	131835	20.000	ng/ul	0.00
20) Naphthalene-d8	10.810	136	589728	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	396380	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	892076	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	746484	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	754180	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	15163	4.460	ng/uL	0.00
4) Pyridine-d5	3.828	84	26996m	3.099	ng/ul	0.03
7) Phenol-d5	7.164	99	62635	5.152	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.316	67	246484	31.597	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	197997	22.531	ng/ul	0.00
15) 4-Methylphenol-d8	8.711	113	131191	13.377	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	137751	28.935	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143	144592	26.524	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	222215	22.382	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	179381	13.717	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	1035842	32.230	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	1101873	30.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.940	143	16936m	3.488	ng/ul	0.08
60) Fluorene-d10	15.634	176	859831	32.081	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	129722	19.763	ng/ul	0.00
73) Anthracene-d10	17.498	188	1467055	33.981	ng/ul	0.00
81) Pyrene-d10	19.728	212	1697690	33.858	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1387034	34.616	ng/ul	0.00

Target Compounds Qvalue

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

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