

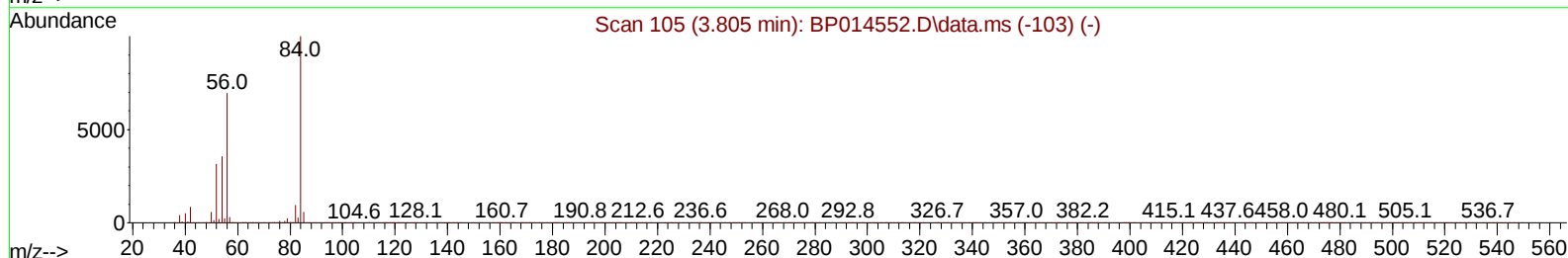
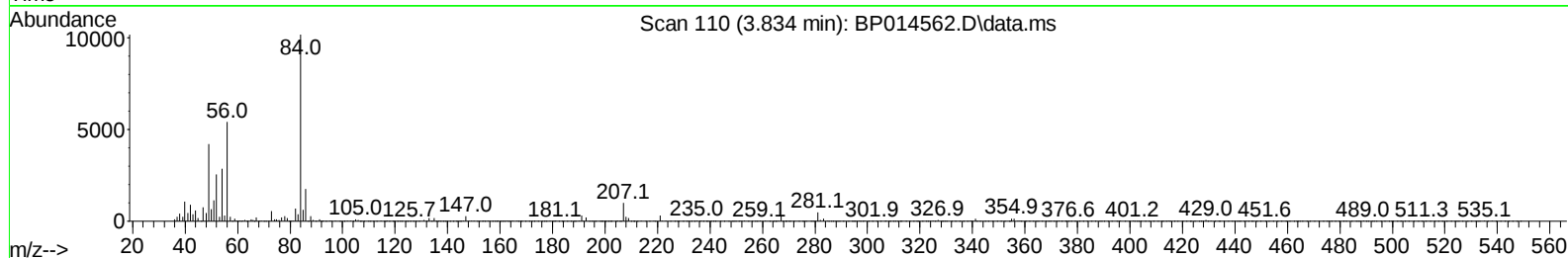
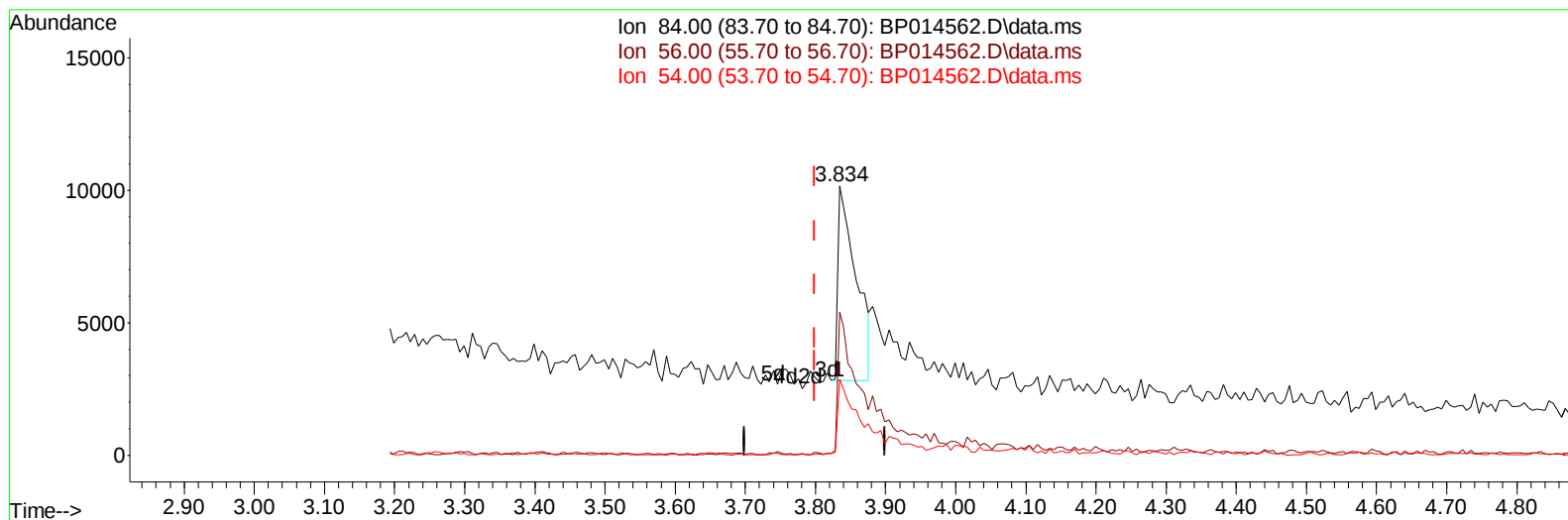
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014562.D
Acq On : 05 Apr 2023 17:44
Operator : CG/JU
Sample : 02122-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB9A8

Manual IntegrationsAPPROVED

Quant Time: Apr 06 00:58:21 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: BP014562.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.035) 1.98 ng/u1

response 13162

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	53.18
54.00	32.60	28.30
0.00	0.00	0.00

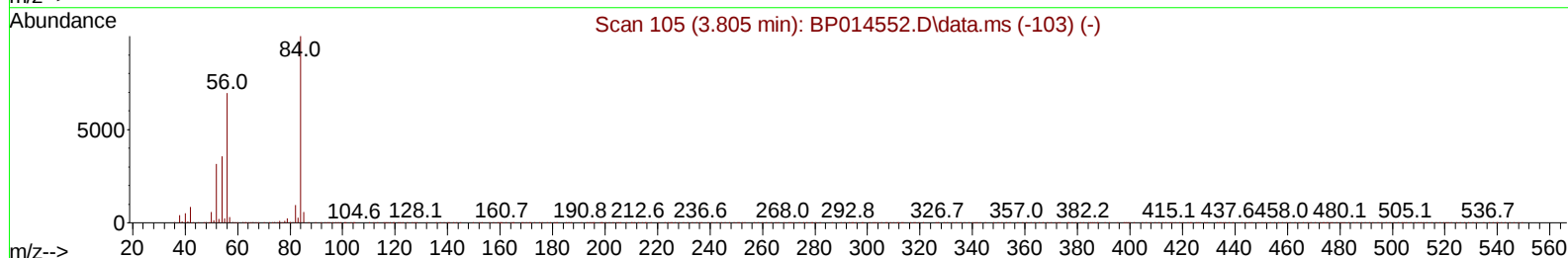
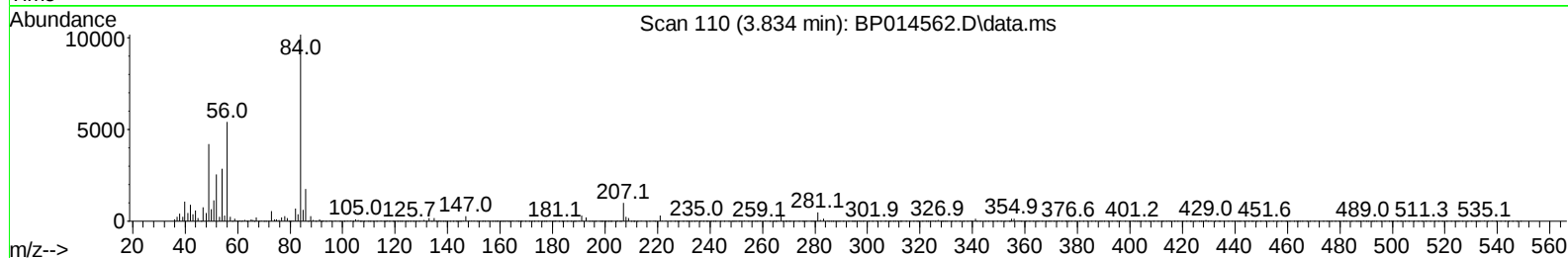
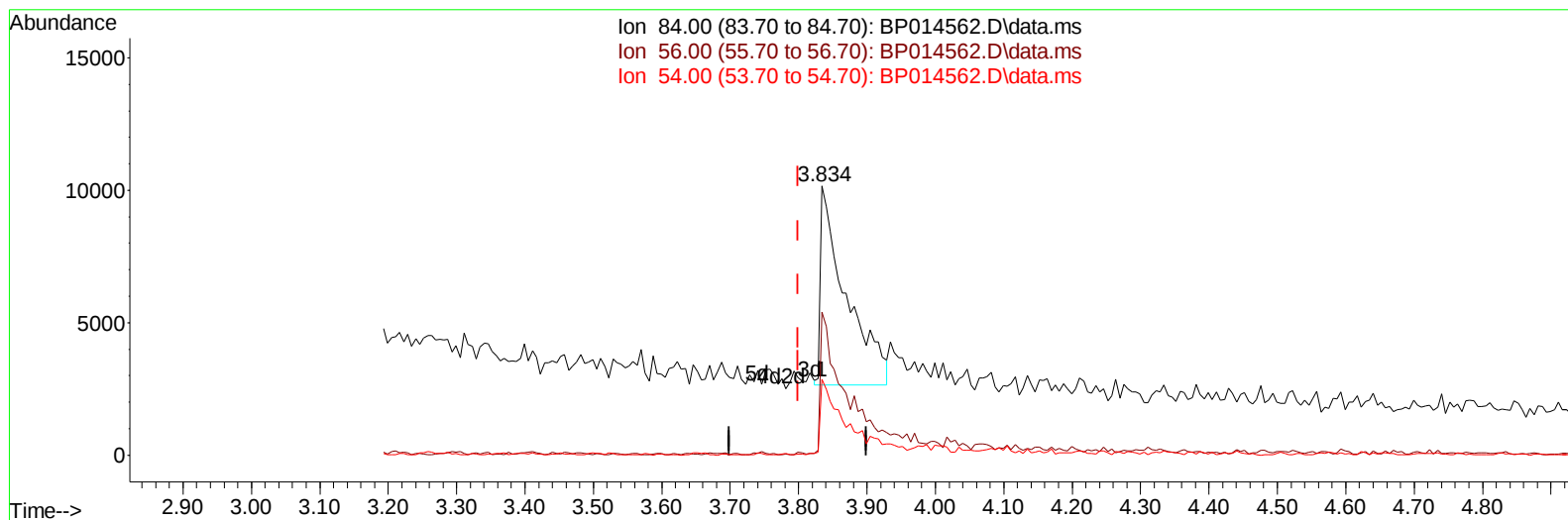
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TIC: BP014562.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.035) 2.92 ng/ul m

response 19371

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	53.18
54.00	32.60	28.30
0.00	0.00	0.00

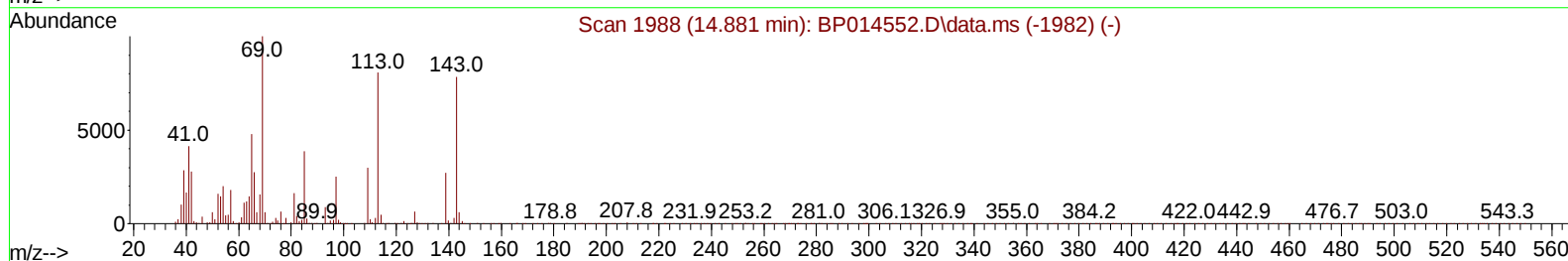
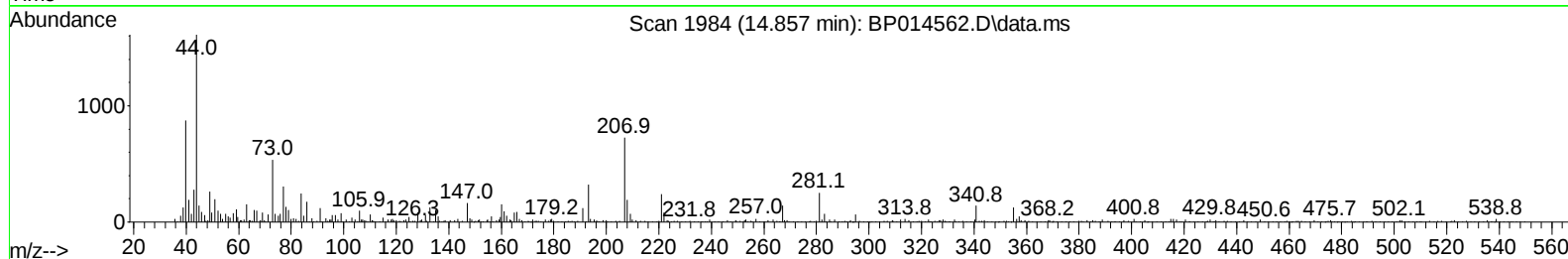
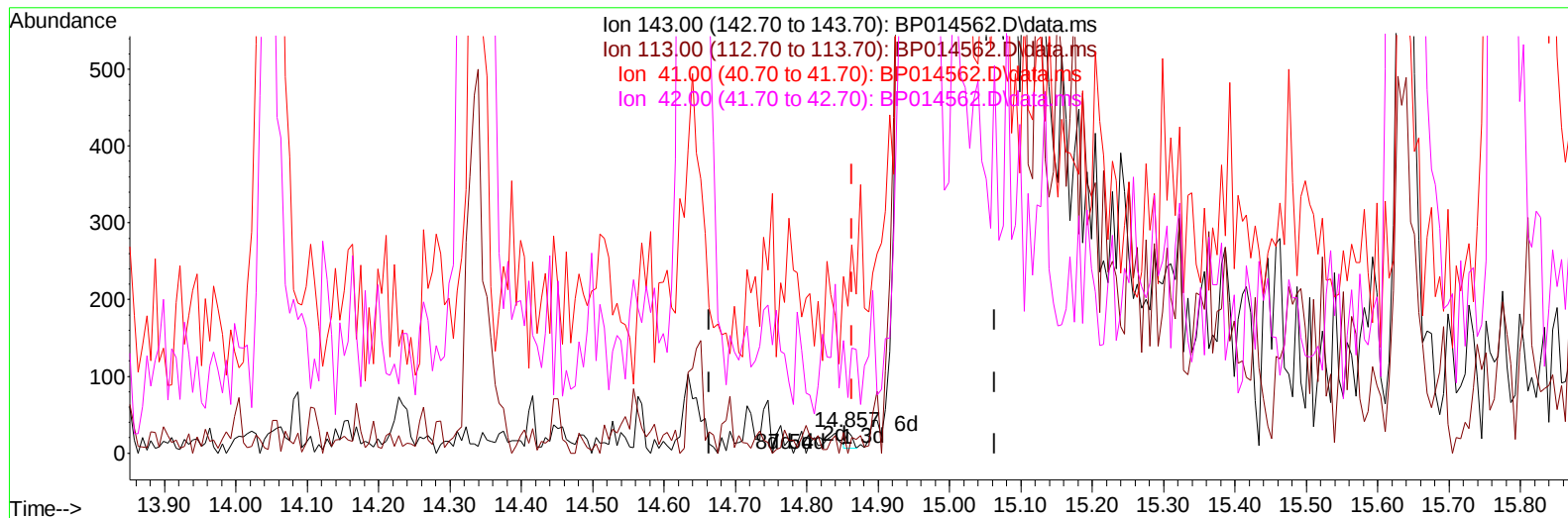
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Misc :
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Instrument :
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ClientSampleId :
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TIC: BP014562.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.006) 0.00 ng/u1

response 13

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	28.57#
41.00	44.20	682.14#
42.00	31.50	257.14#

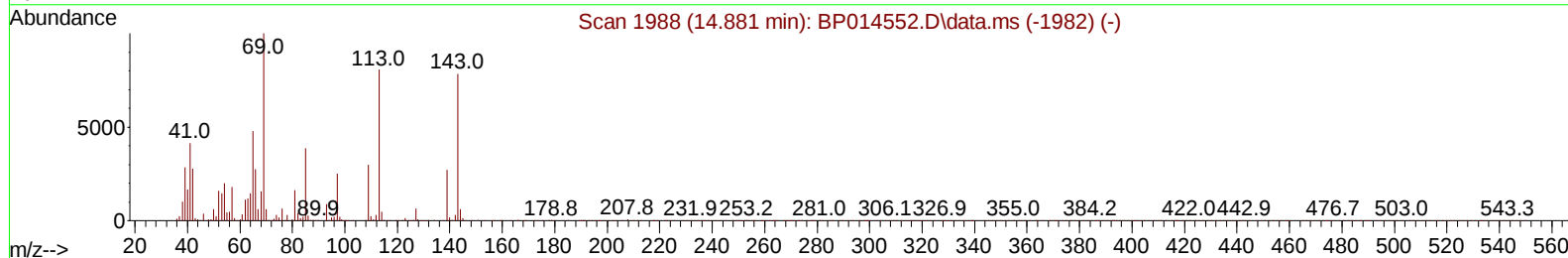
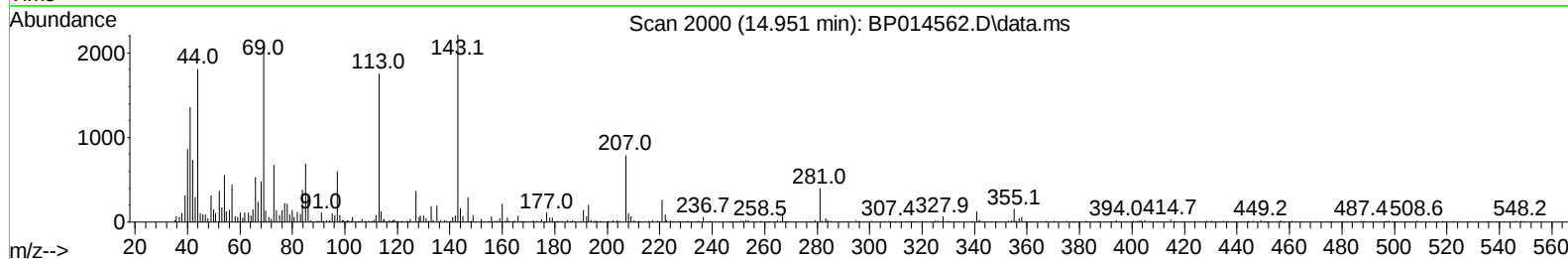
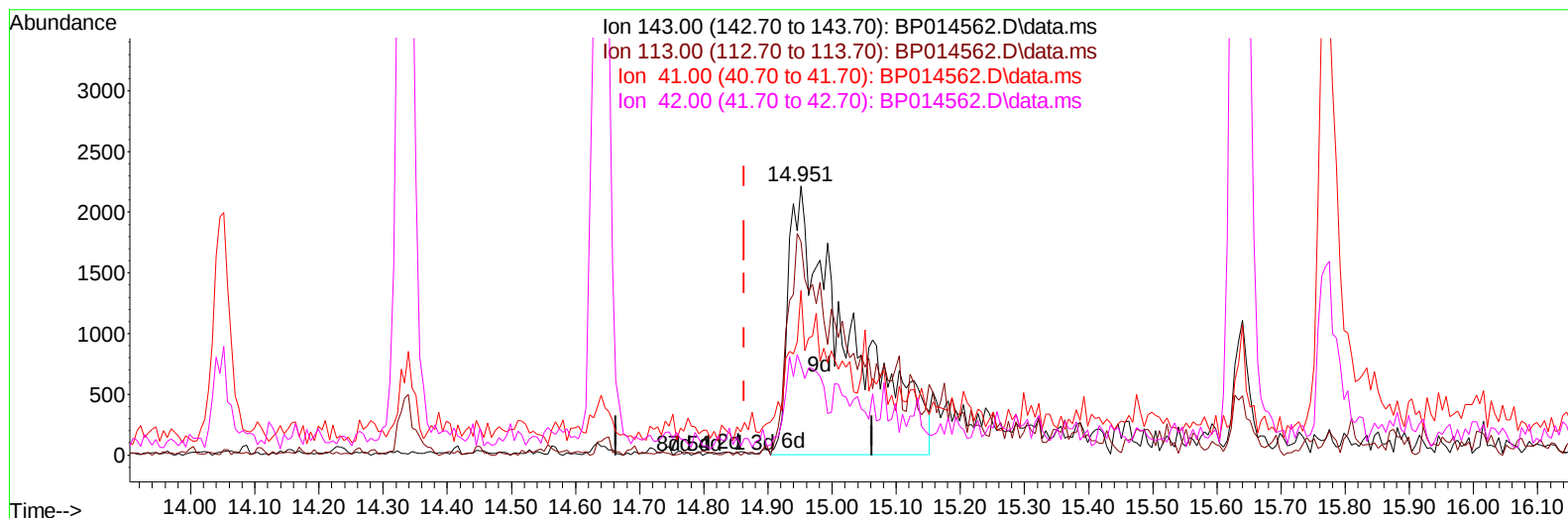
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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TIC: BP014562.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.951min (+ 0.088) 4.05 ng/ul m

response 14183

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	79.37
41.00	44.20	61.26#
42.00	31.50	33.27

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 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	100397	20.000	ng/ul	0.00
20) Naphthalene-d8	10.810	136	432126	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	285603	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	605438	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	550942	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	631370	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	11796	4.556	ng/uL	0.00
4) Pyridine-d5	3.834	84	19371m	2.920	ng/ul	0.04
7) Phenol-d5	7.169	99	53852	5.816	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.311	67	214207	36.058	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	174945	26.142	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	105258	14.093	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	116937	33.521	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143	116406	29.141	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	186926	25.695	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	110965	11.580	ng/ul	0.01
46) Dimethylphthalate-d6	14.051	166	829343	35.814	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	886483	34.500	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	14183m	4.054	ng/ul	0.09
60) Fluorene-d10	15.634	176	690391	35.750	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	96780	21.724	ng/ul	0.00
73) Anthracene-d10	17.504	188	1131609	38.621	ng/ul	0.00
81) Pyrene-d10	19.733	212	1330803	35.961	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1287107	38.370	ng/ul	0.00

Target Compounds Qvalue

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

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