

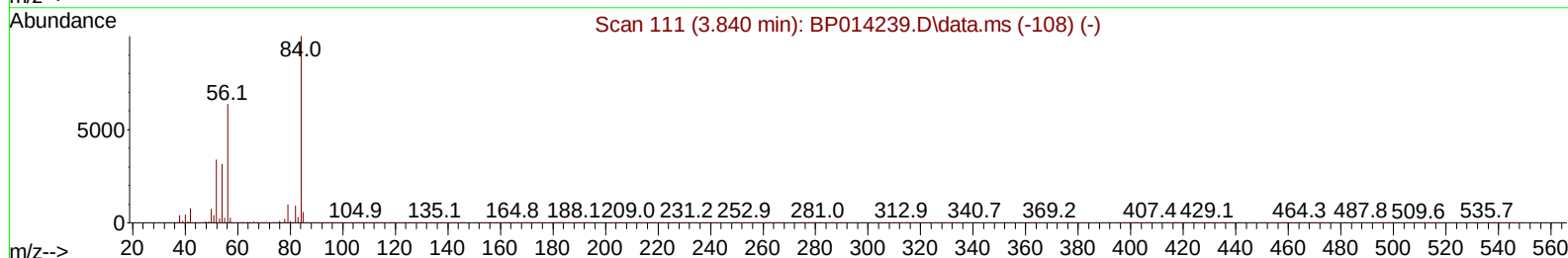
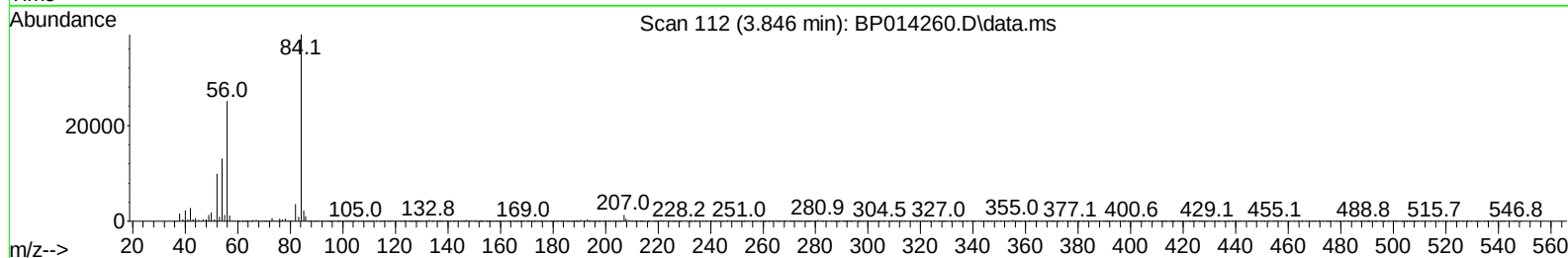
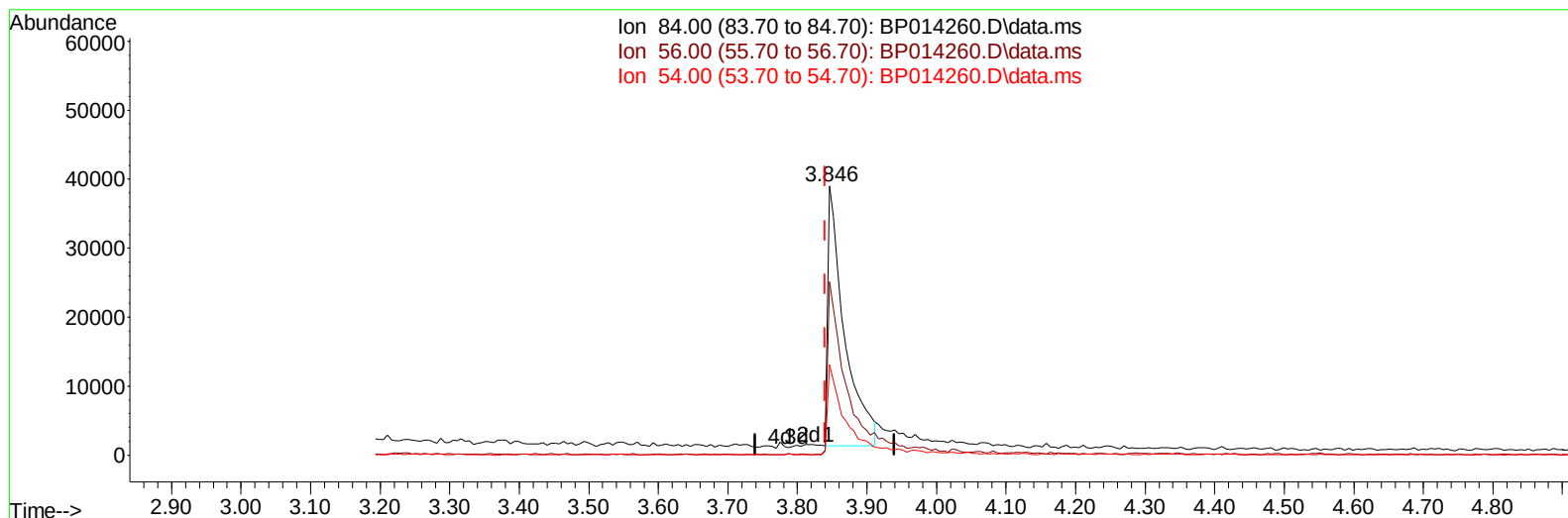
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014260.D
 Acq On : 23 Mar 2023 16:25
 Operator : CG/JU
 Sample : 02010-02
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB945

Manual IntegrationsAPPROVED

Quant Time: Mar 24 00:51:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 23 04:56:06 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/24/2023
 Supervised By :Jagrut Upadhyay 03/24/2023



TIC: BP014260.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 6.43 ng/ul

response 62287

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	64.39
54.00	33.20	33.64
0.00	0.00	0.00

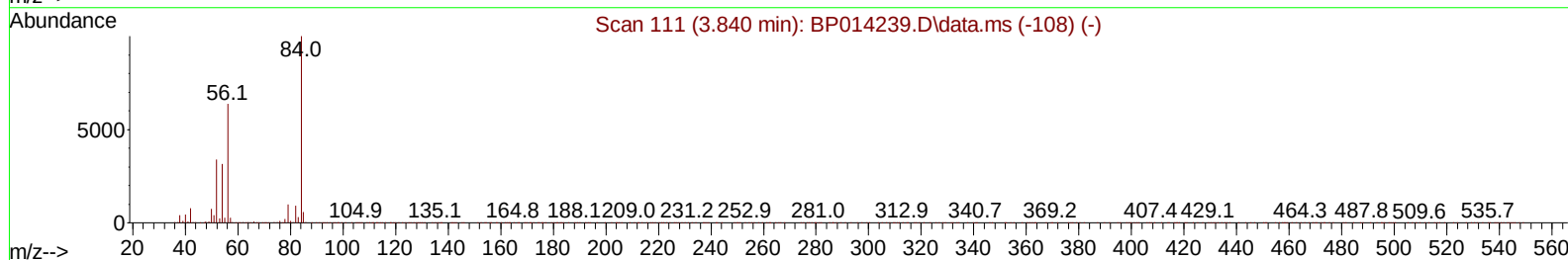
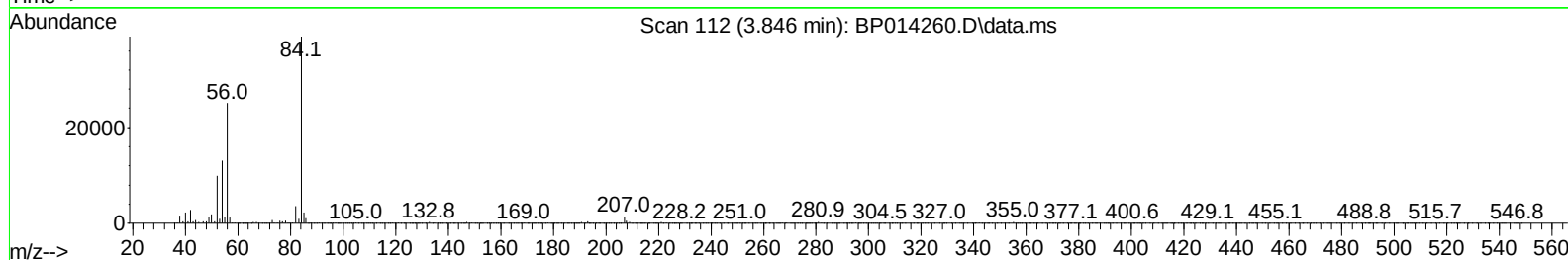
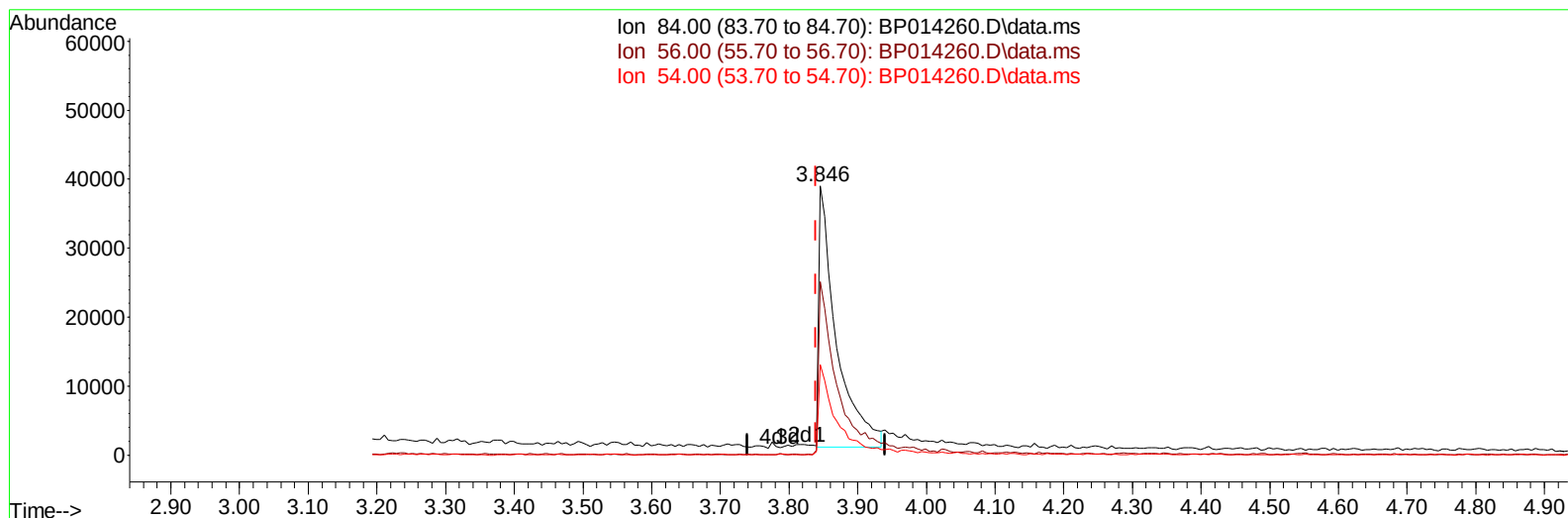
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TIC: BP014260.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 6.86 ng/ul m

response 66492

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	64.39
54.00	33.20	33.64
0.00	0.00	0.00

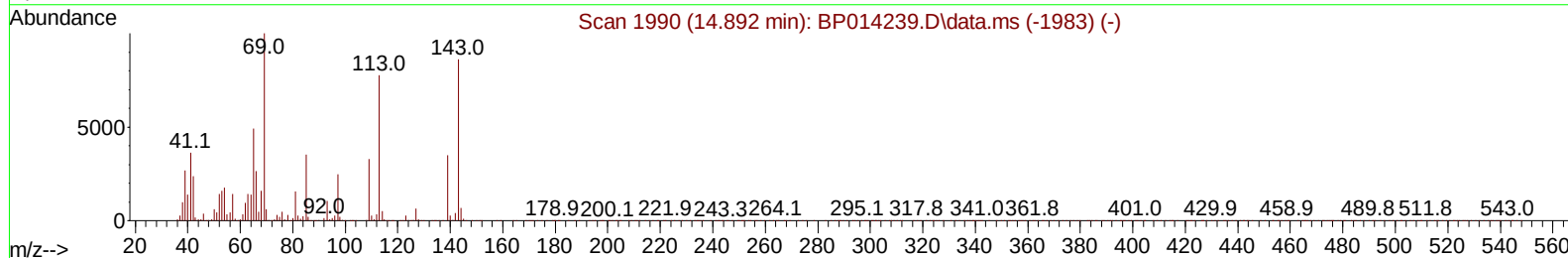
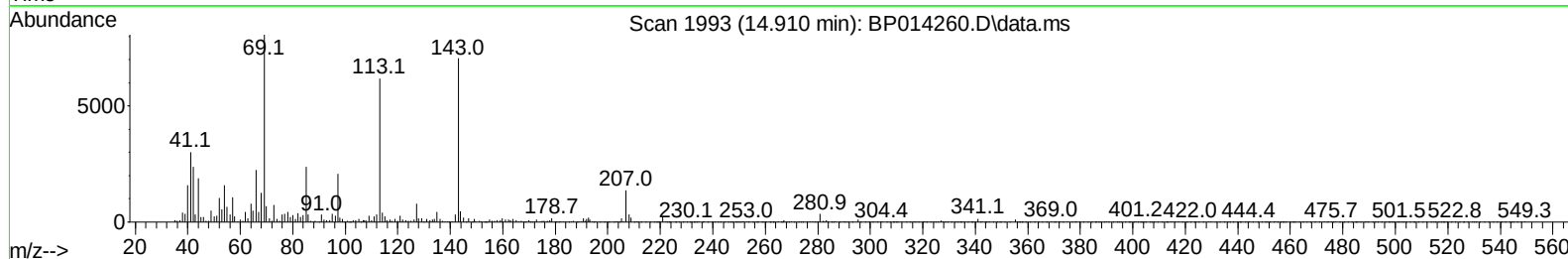
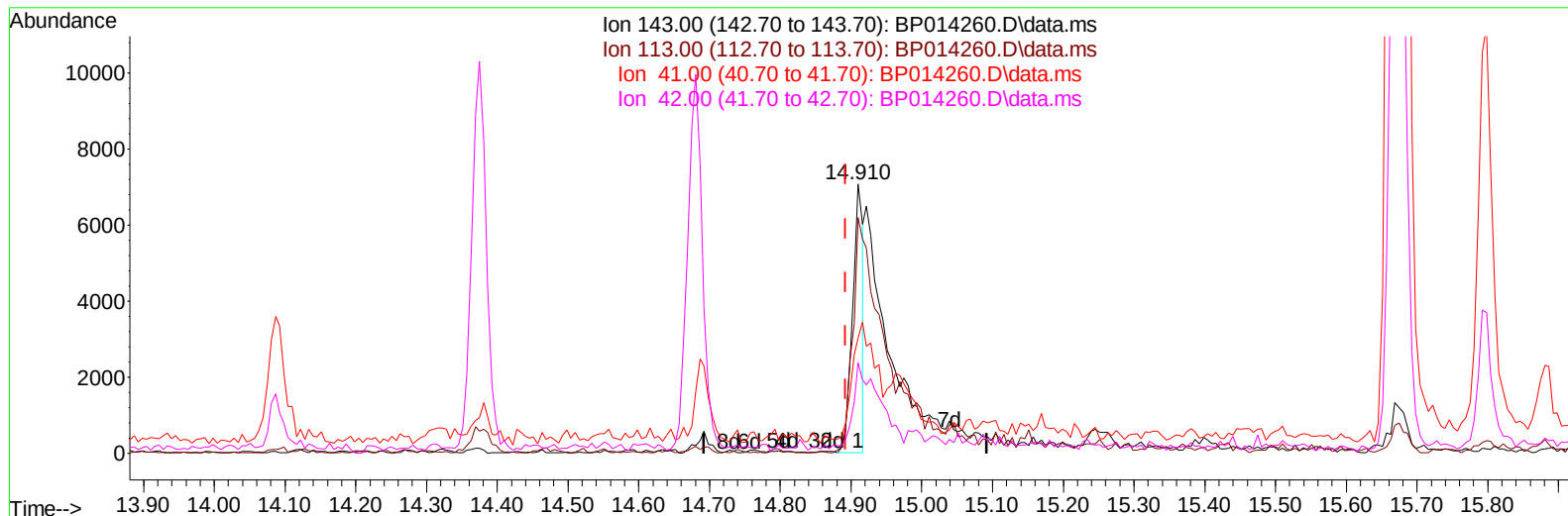
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TIC: BP014260.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.018) 1.09 ng/ul

response 7292

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	87.50
41.00	40.20	42.76
42.00	28.50	33.52

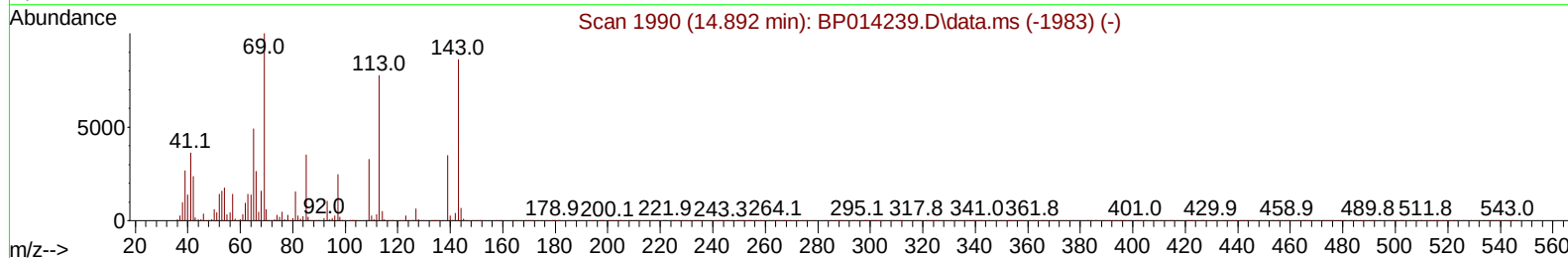
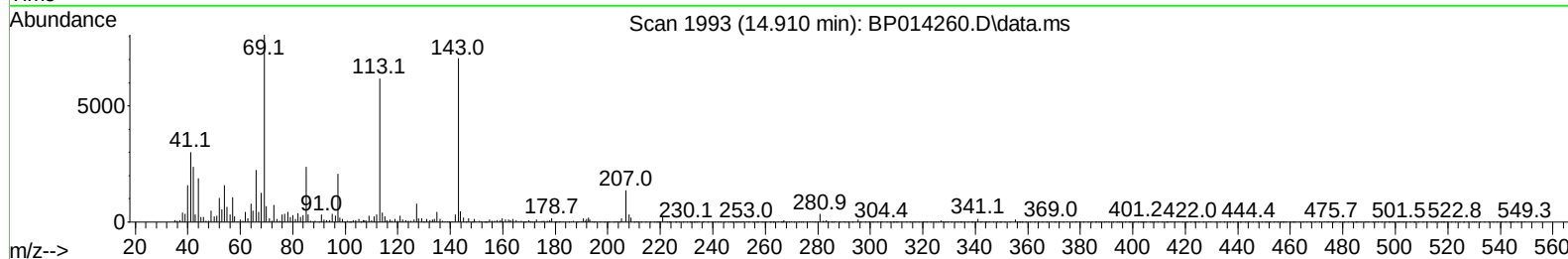
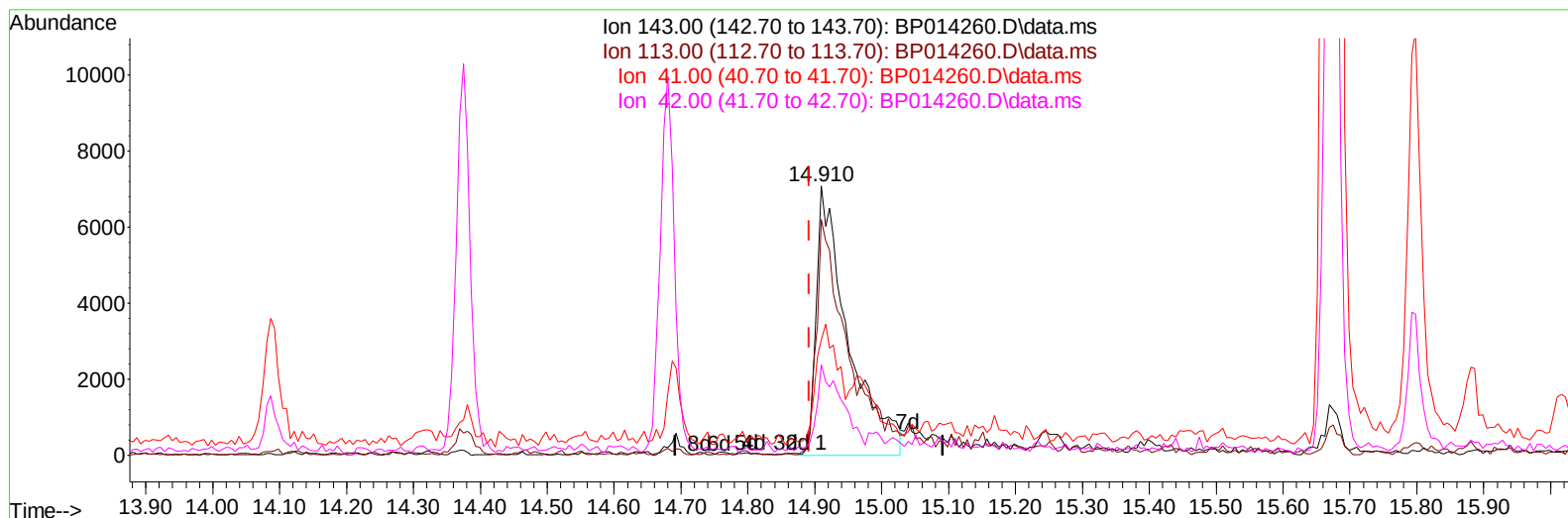
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014260.D
Acq On : 23 Mar 2023 16:25
Operator : CG/JU
Sample : 02010-02
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP014260.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.910min (+ 0.018) 3.46 ng/ul m

response 23160

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	87.50
41.00	40.20	42.76
42.00	28.50	33.52

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 Sample : 02010-02
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	132233	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	610996	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.681	164	447345	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.439	188	1067090	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.521	240	1095173	20.000	ng/ul	# 0.00
88) Perylene-d12	24.039	264	1142250	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	13149	3.757	ng/uL	0.00
4) Pyridine-d5	3.846	84	66492m	6.862	ng/ul	0.00
7) Phenol-d5	7.199	99	56817	4.869	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.357	67	207814	30.536	ng/ul	0.00
11) 2-Chlorophenol-d4	7.563	132	172324	19.383	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	121300	12.726	ng/ul	0.00
21) Nitrobenzene-d5	9.210	128	127262	26.104	ng/ul	0.00
24) 2-Nitrophenol-d4	9.934	143	138580	23.747	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	232418	21.166	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	313265	21.558	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	1111941	29.289	ng/ul	0.00
49) Acenaphthylene-d8	14.375	160	1130984	28.122	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	23160m	3.462	ng/ul	0.02
60) Fluorene-d10	15.675	176	936797	29.115	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	188026	23.313	ng/ul	0.00
73) Anthracene-d10	17.539	188	1642997	31.717	ng/ul	0.00
81) Pyrene-d10	19.763	212	2031347	33.630	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	1964349	32.330	ng/ul	0.00

Target Compounds Qvalue

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

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