

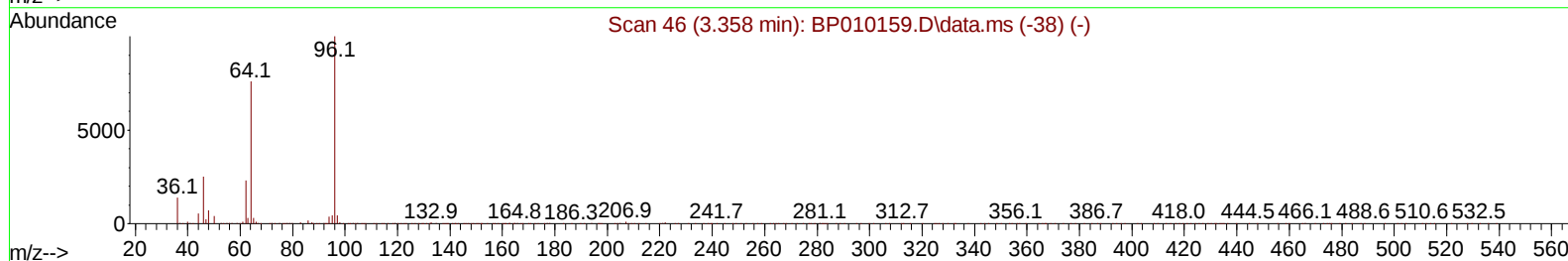
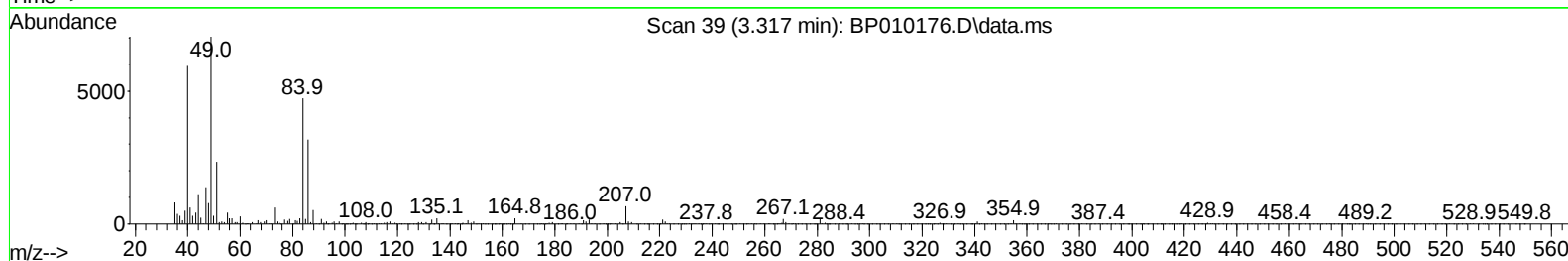
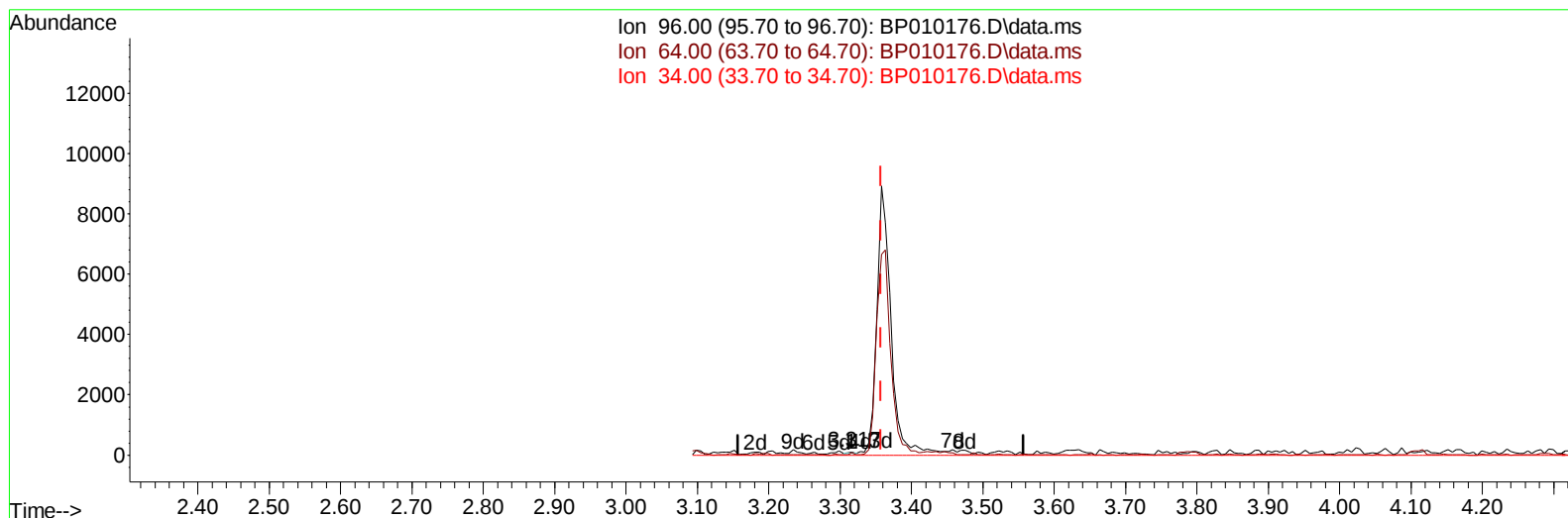
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010176.D
 Acq On : 02 May 2022 16:58
 Operator : CG/JU
 Sample : N2562-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW464

Manual IntegrationsAPPROVED

Quant Time: May 03 01:37:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/03/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010176.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.317min (-0.041) 0.01 ng/uL

response 26

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	65.42
34.00	0.00	0.00
0.00	0.00	0.00

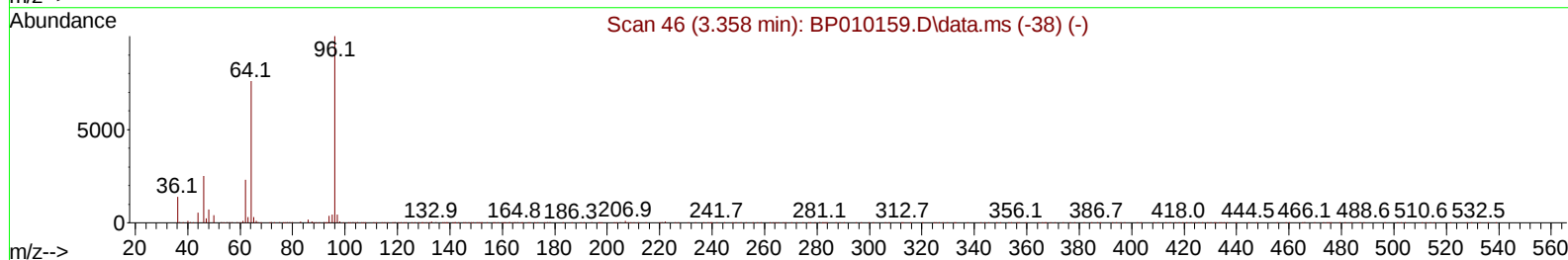
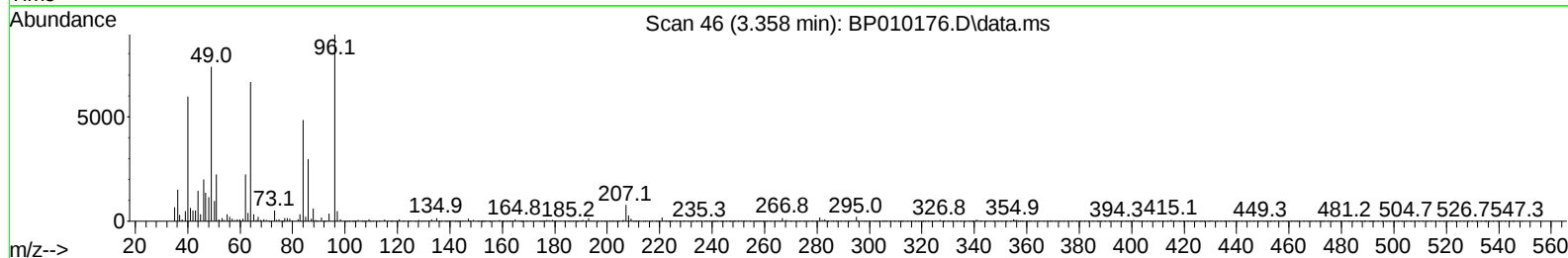
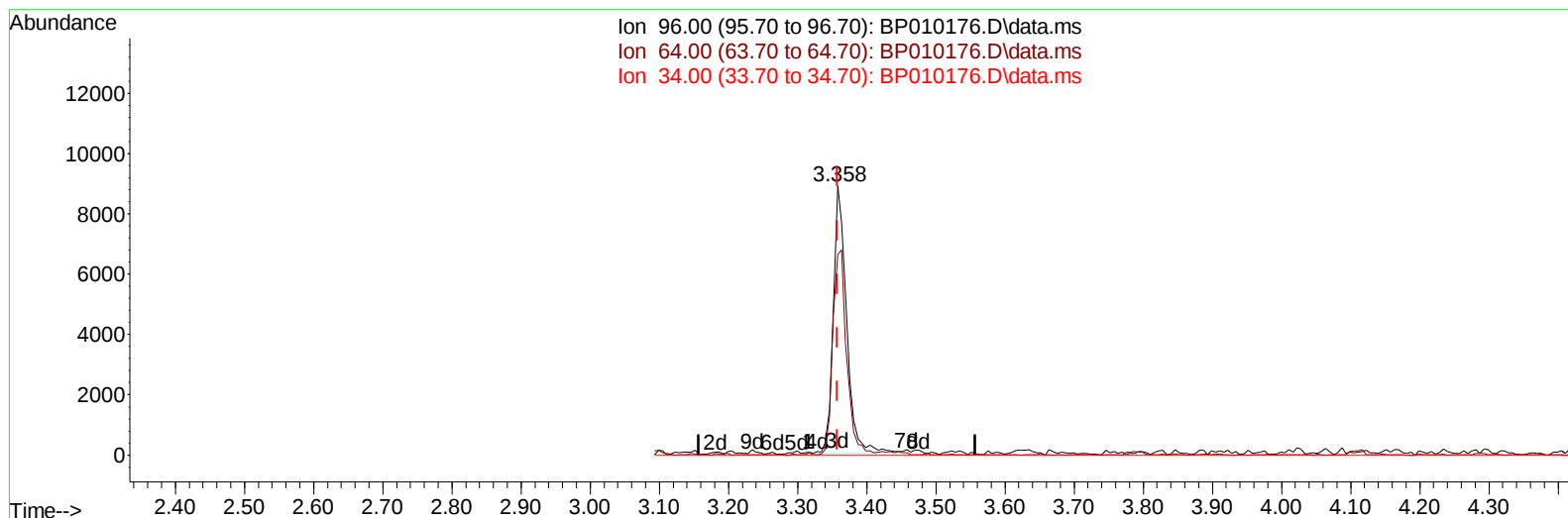
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TIC: BP010176.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.358min (-0.000) 2.94 ng/uL m

response 11789

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.54#
34.00	0.00	0.00
0.00	0.00	0.00

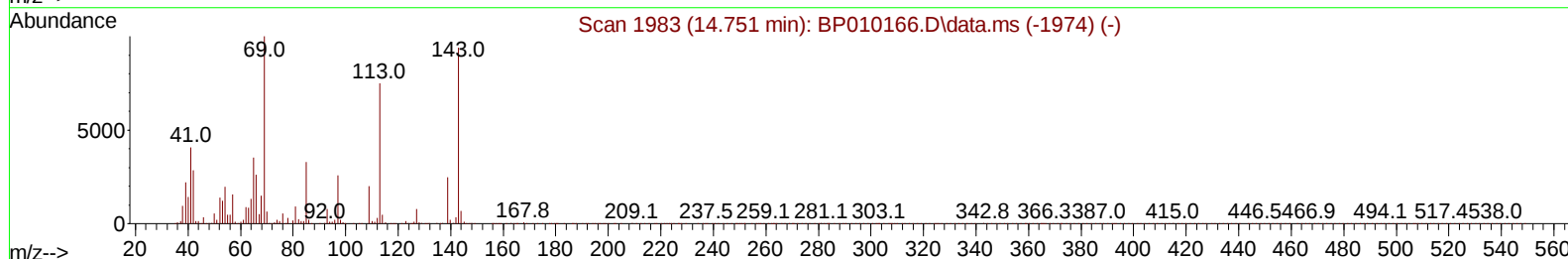
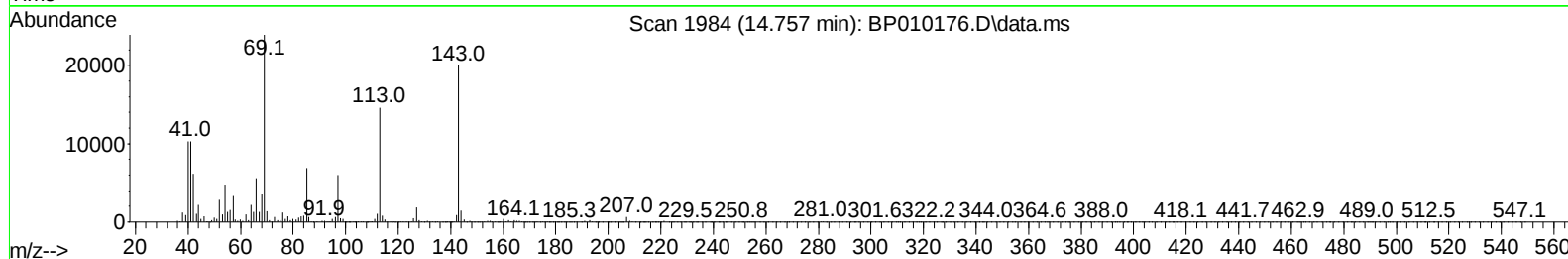
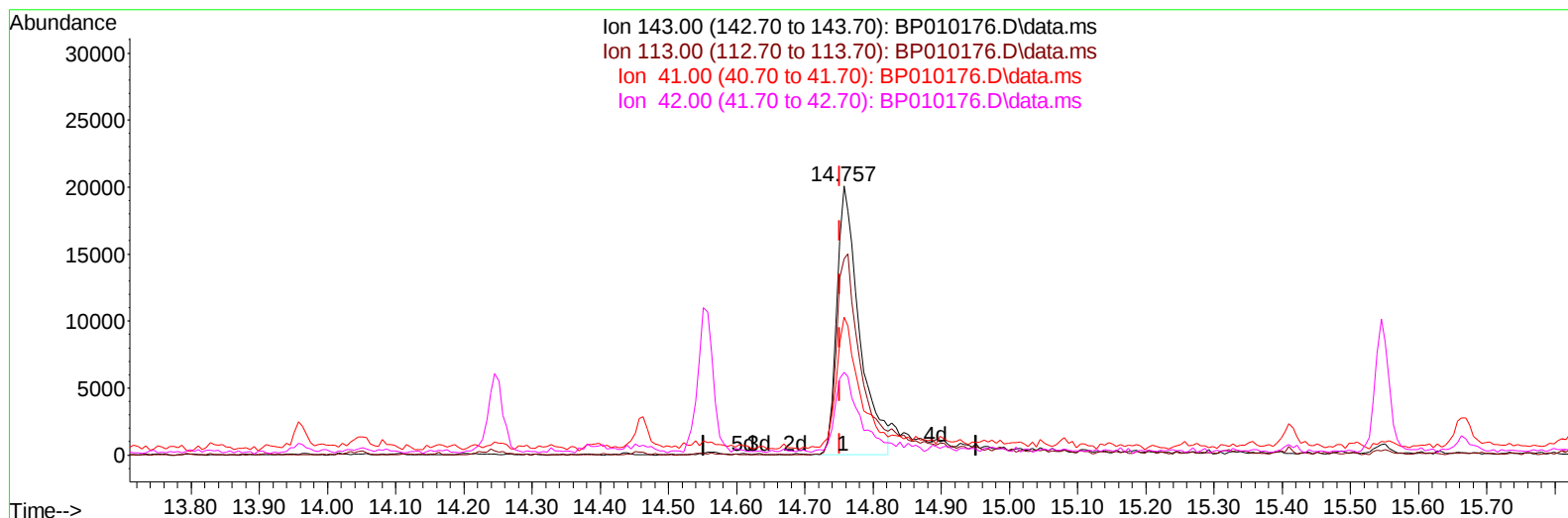
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TIC: BP010176.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 6.61 ng/ul

response 46446

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.87#
41.00	23.20	51.31#
42.00	18.00	30.81#

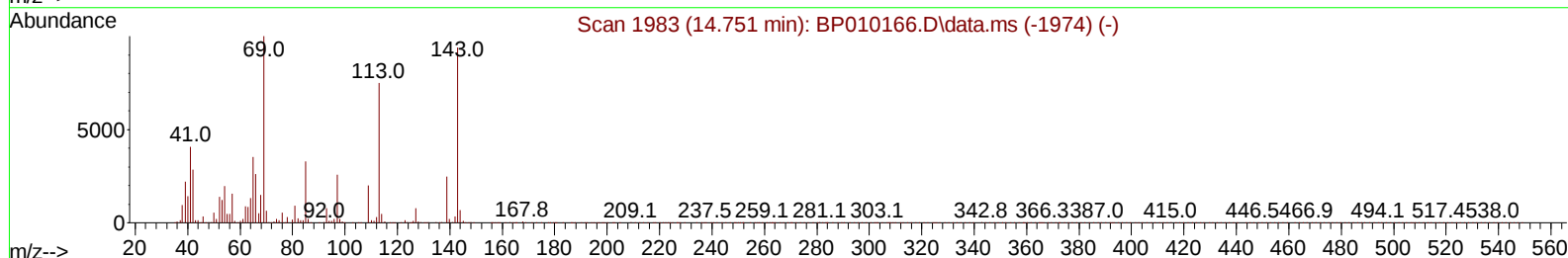
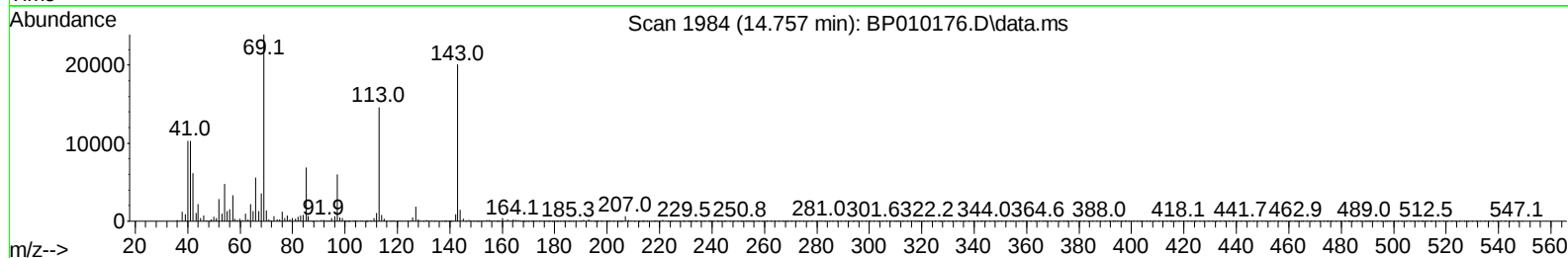
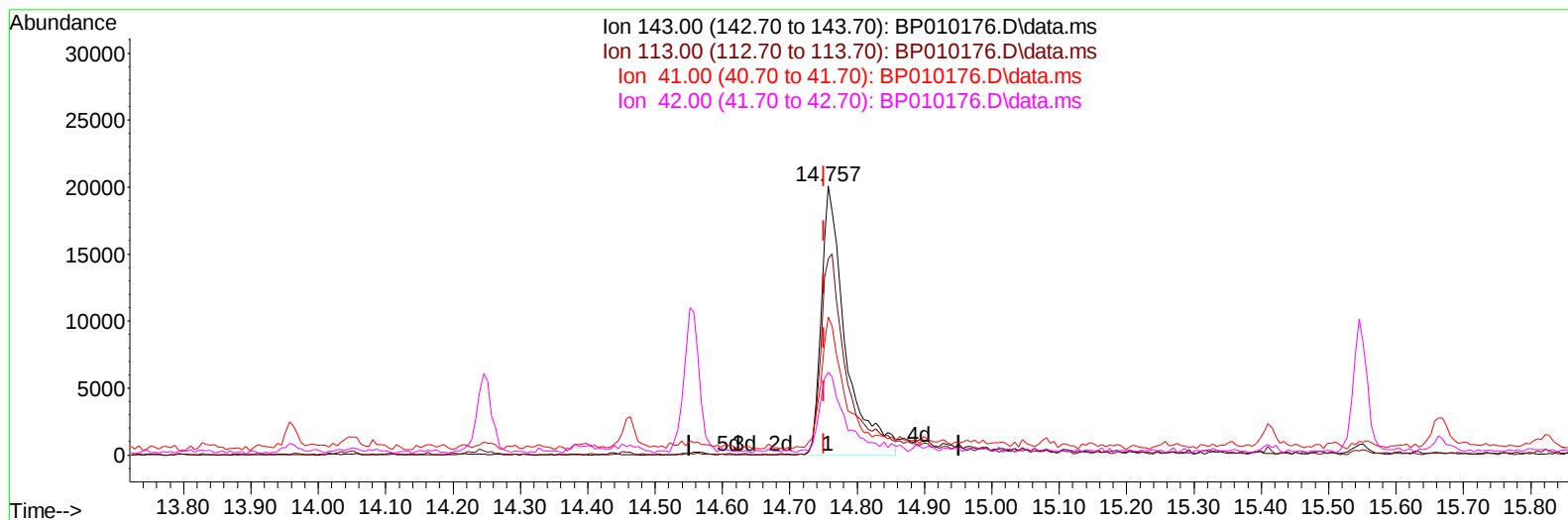
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 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW464

Manual IntegrationsAPPROVED

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TIC: BP010176.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (+ 0.006) 7.15 ng/ul m

response 50185

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.87#
41.00	23.20	51.31#
42.00	18.00	30.81#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	174563	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	752734	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	490351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	928264	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	686376	20.000	ng/ul	# 0.00
88) Perylene-d12	23.833	264	607952	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	11789m	2.945	ng/uL	0.00
4) Pyridine-d5	3.787	84	55549	4.933	ng/ul	0.01
7) Phenol-d5	7.087	99	213626	13.454	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	156696	16.447	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	185125	15.743	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	170358	12.950	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	96603	16.136	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	105720	16.169	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	173068	13.597	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	231894	12.706	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	611397	16.029	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	695311	15.062	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	50185m	7.147	ng/ul	0.00
60) Fluorene-d10	15.545	176	525430	16.014	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	43513	7.413	ng/ul	0.00
73) Anthracene-d10	17.404	188	718716	16.004	ng/ul	0.00
81) Pyrene-d10	19.639	212	796973	20.417	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	565105	17.730	ng/ul	0.00

Target Compounds Qvalue

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

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