

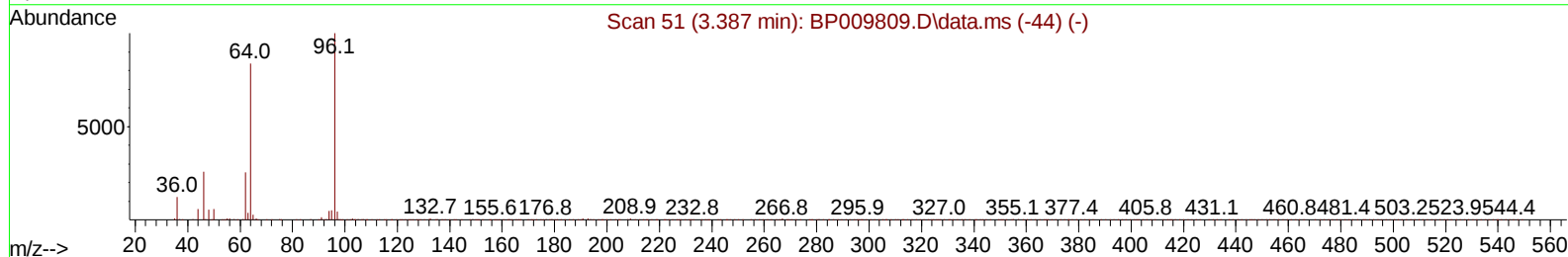
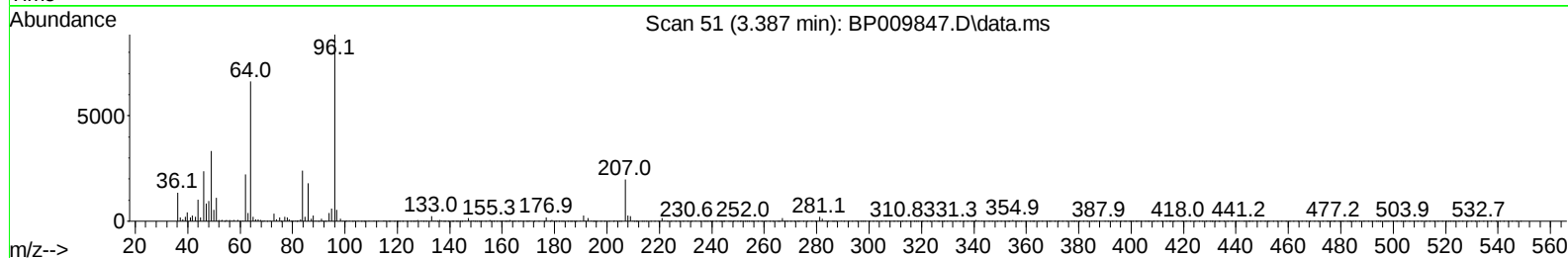
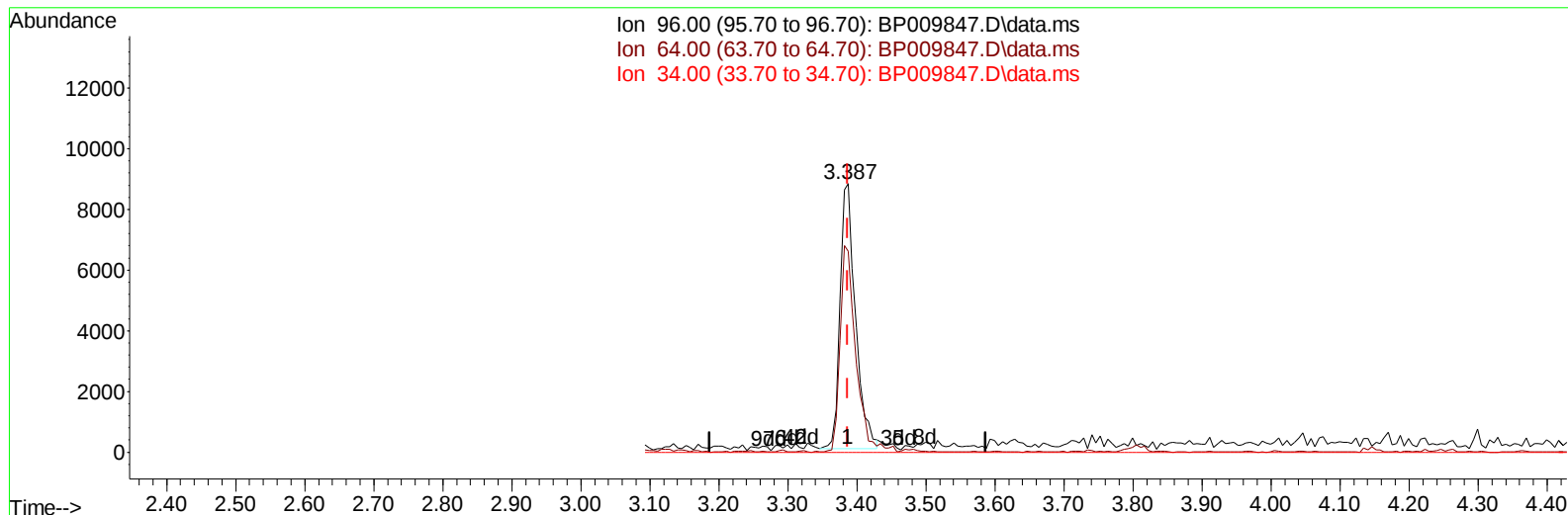
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
Data File : BP009847.D
Acq On : 15 Apr 2022 02:32
Operator : CG/JU
Sample : N2293-18
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ETNP6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/15/2022
Supervised By :Yogesh Patel 04/22/2022

Quant Time: Apr 15 04:22:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration



TIC: BP009847.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.387min (+ 0.000) 4.06 ng/uL****response 13639**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.82#
34.00	0.00	0.00
0.00	0.00	0.00

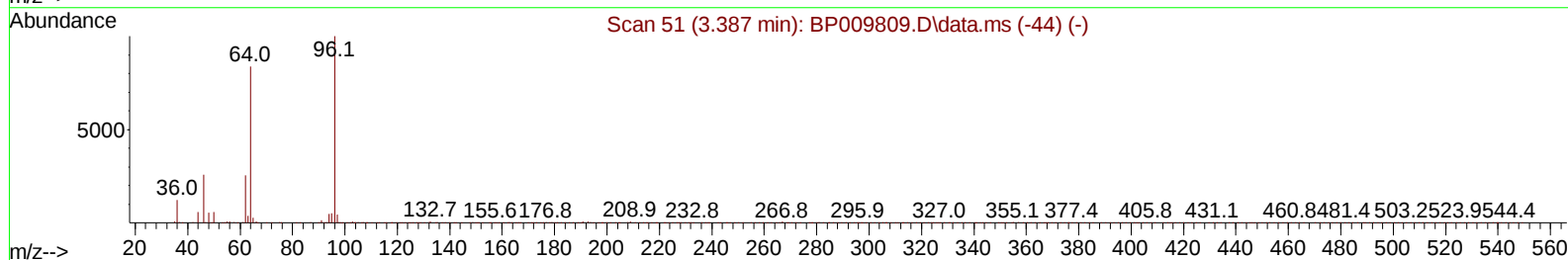
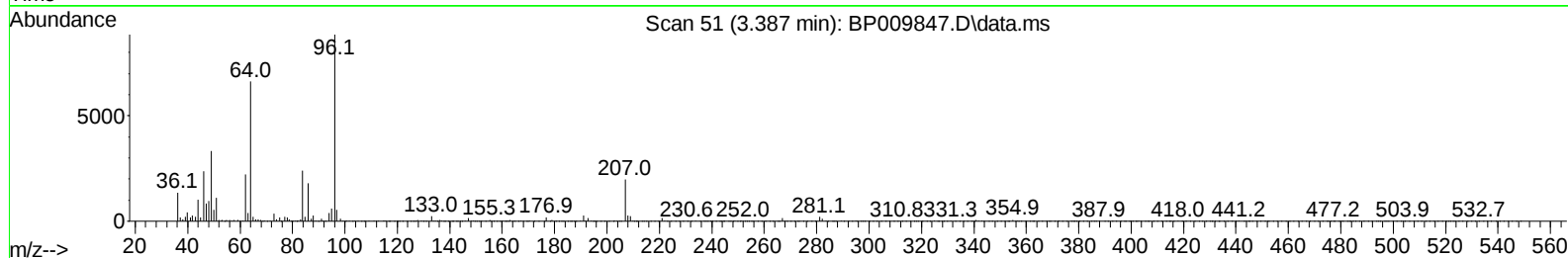
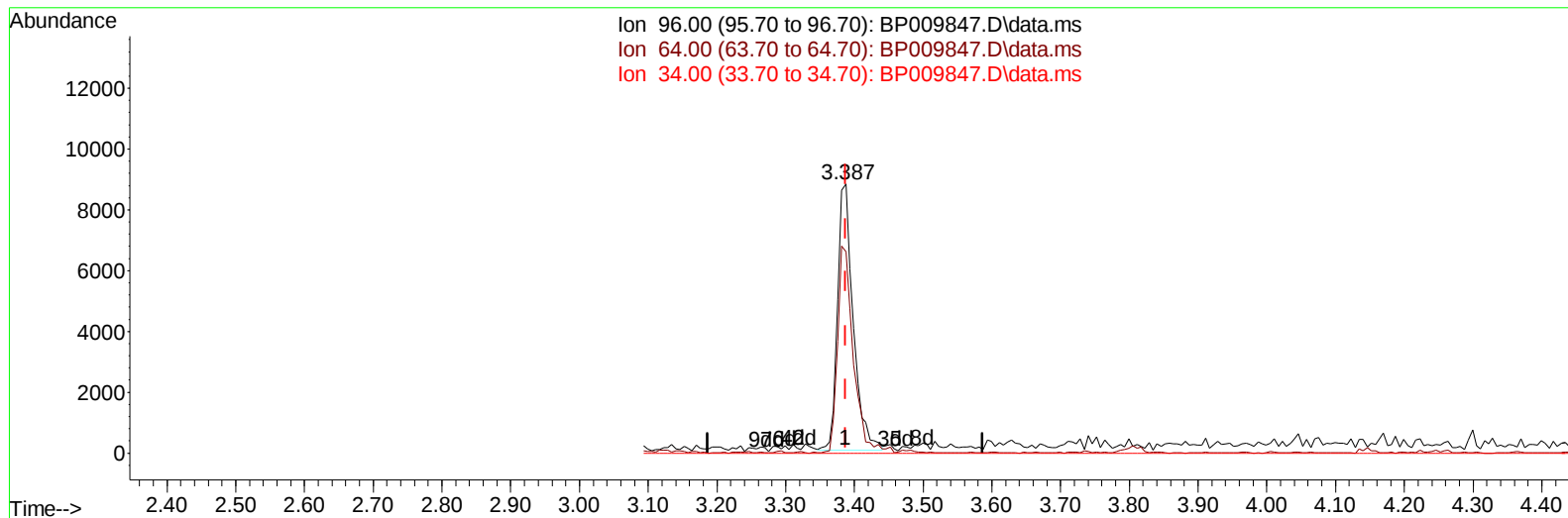
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TIC: BP009847.D\data.ms

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

3.387min (+ 0.000) 4.12 ng/uL m

response 13868

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	74.82#
34.00	0.00	0.00
0.00	0.00	0.00

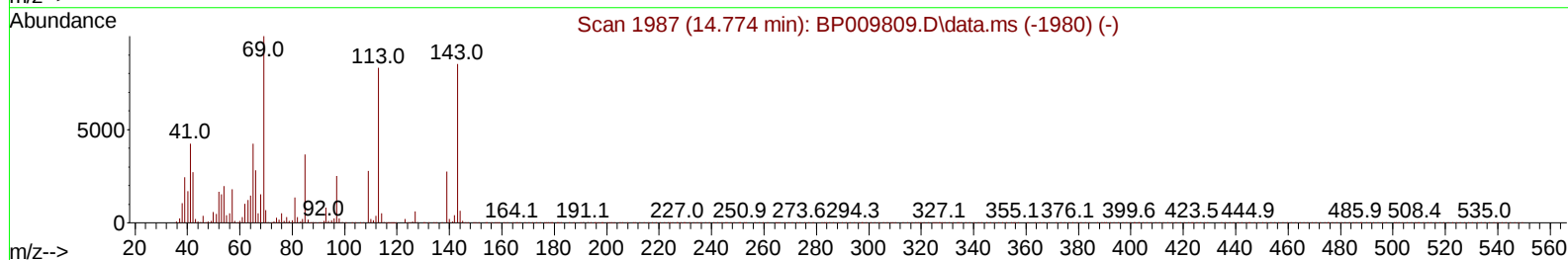
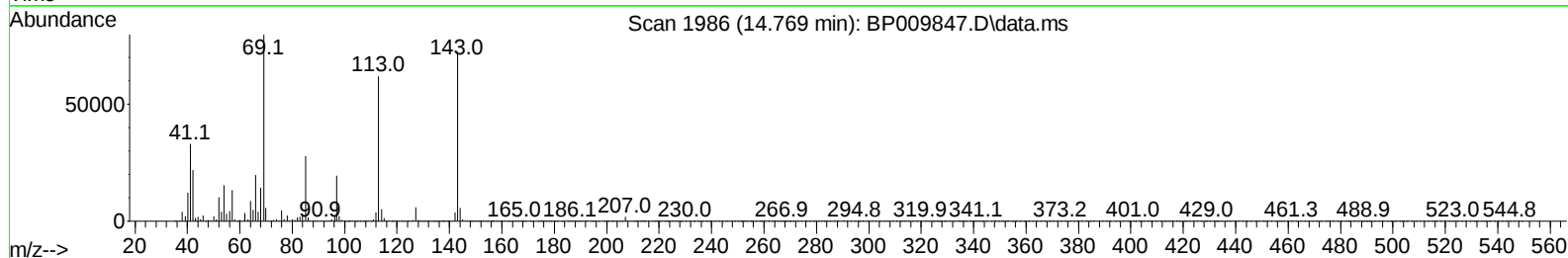
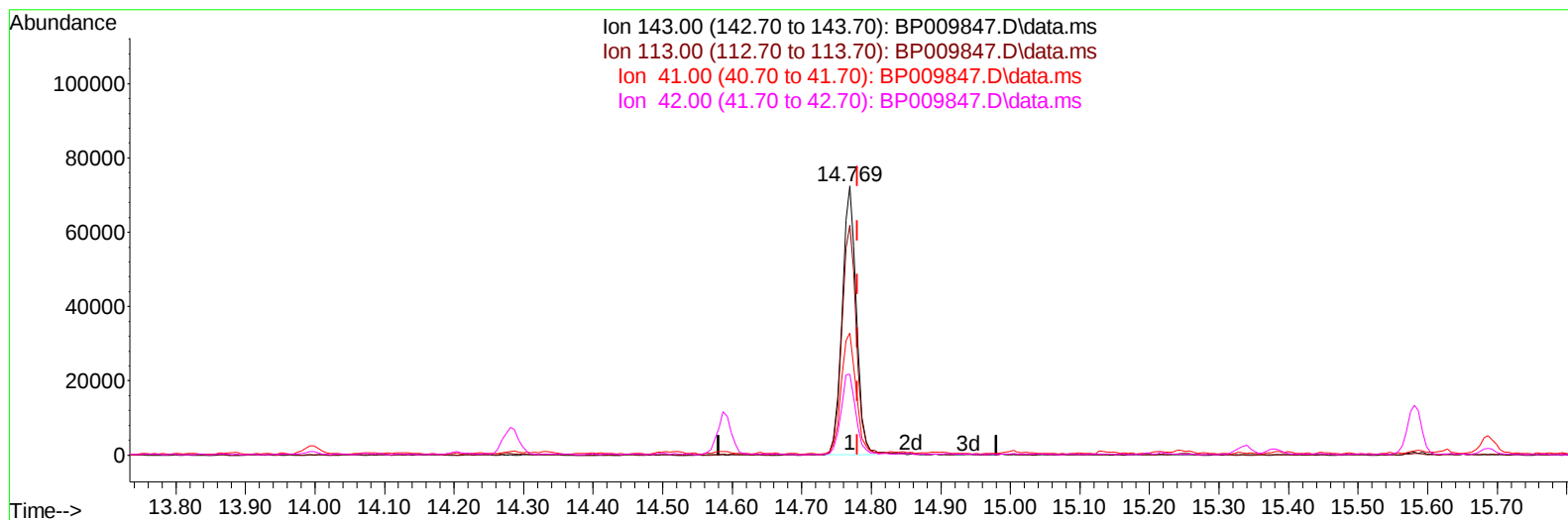
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TIC: BP009847.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.769min (-0.012) 15.36 ng/u1****response 102263**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.37#
41.00	23.20	45.33#
42.00	18.00	30.07#

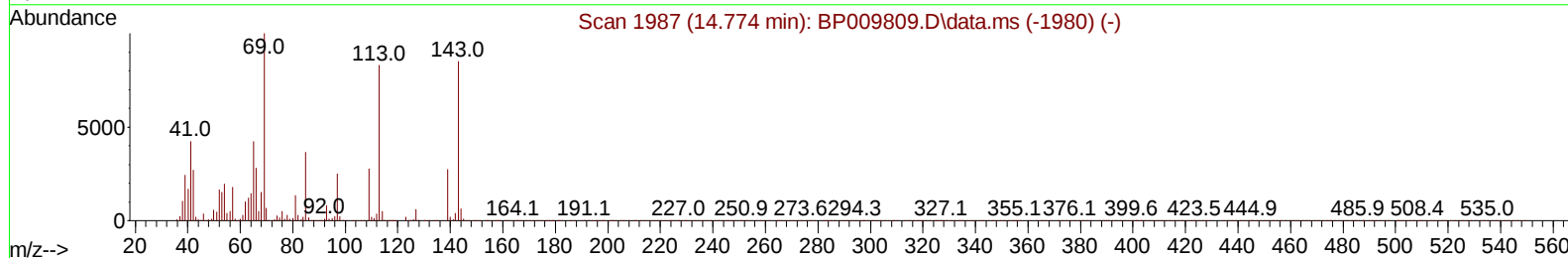
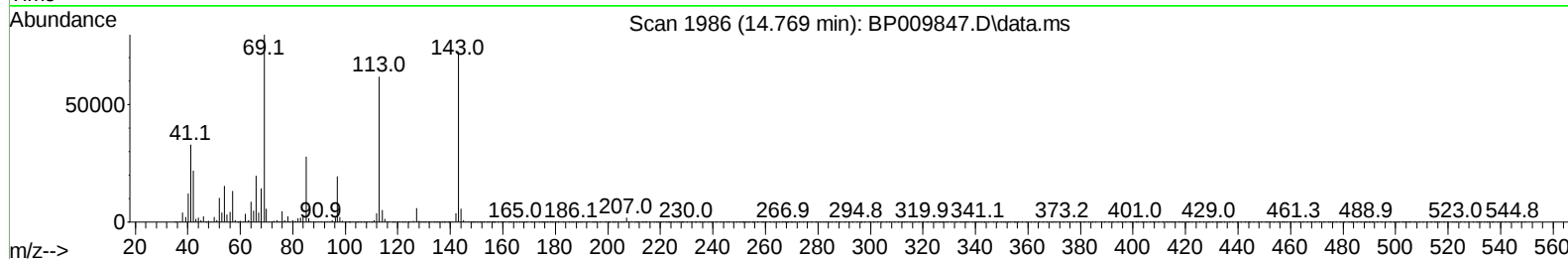
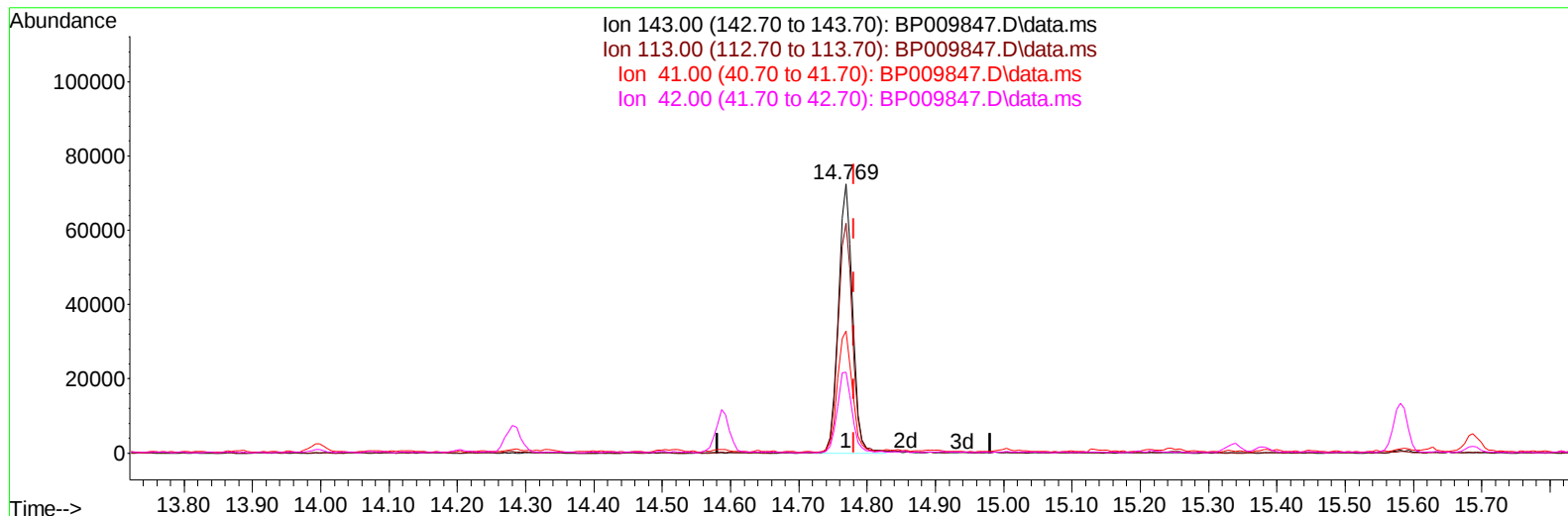
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 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
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TIC: BP009847.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (-0.012) 15.61 ng/ul m

response 103938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.37#
41.00	23.20	45.33#
42.00	18.00	30.07#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.964	152	150094	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.769	136	658692	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	467424	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.340	188	1043835	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.416	240	1036744	20.000	ng/ul	#-0.02
88) Perylene-d12	23.880	264	904868	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	13868m	4.124	ng/uL	0.00
4) Pyridine-d5	3.805	84	110563	11.924	ng/ul	0.00
7) Phenol-d5	7.111	99	267021	20.264	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	171081	21.808	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.487	132	213687	21.498	ng/ul	-0.01
15) 4-Methylphenol-d8	8.658	113	234948	21.358	ng/ul	-0.01
21) Nitrobenzene-d5	9.116	128	112909	23.770	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.840	143	121431	23.318	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	228207	20.941	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.899	131	292754	18.862	ng/ul	0.00
46) Dimethylphthalate-d6	13.999	166	804465	21.910	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	896931	20.620	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	103938m	15.609	ng/ul	-0.01
60) Fluorene-d10	15.581	176	684930	22.175	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.687	200	74860	12.106	ng/ul	-0.01
73) Anthracene-d10	17.440	188	1086381	21.737	ng/ul	-0.01
81) Pyrene-d10	19.663	212	1313040	23.190	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.722	264	1038250	22.033	ng/ul	-0.02
Target Compounds						
43) 1,1'-Biphenyl	13.428	154	53209	1.560	ng/ul	Qvalue 95

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

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