

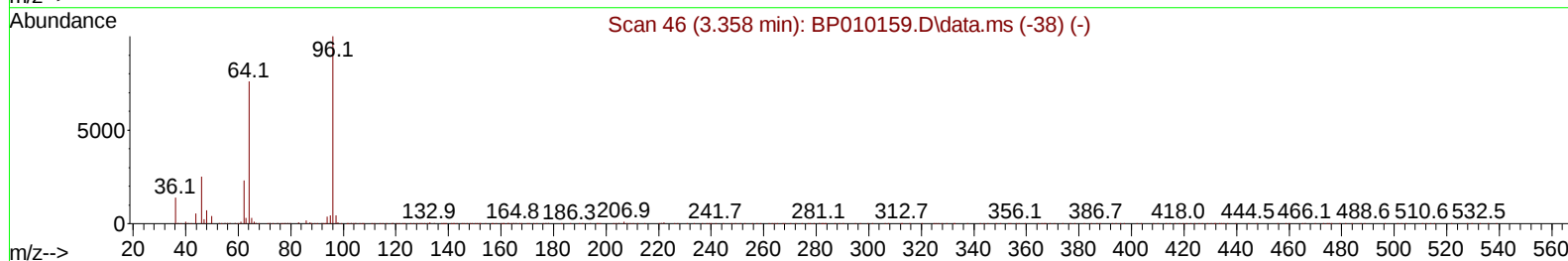
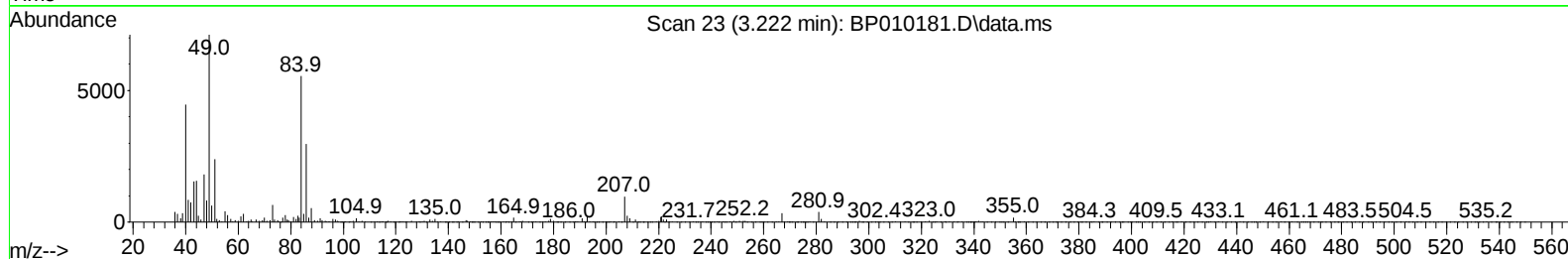
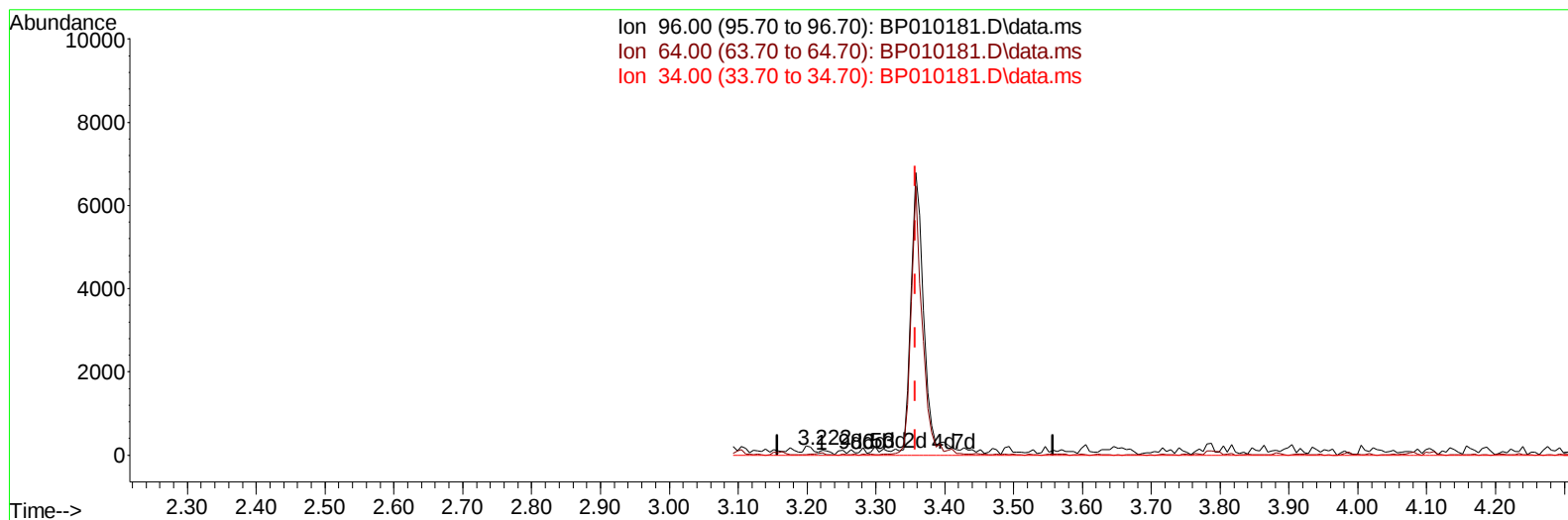
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010181.D
 Acq On : 02 May 2022 19:47
 Operator : CG/JU
 Sample : N2562-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW463

Manual IntegrationsAPPROVED

Quant Time: May 03 01:38:05 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: BP010181.D\data.ms

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

3.222min (-0.135) 0.03 ng/uL

response 90

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.45
34.00	0.00	0.00
0.00	0.00	0.00

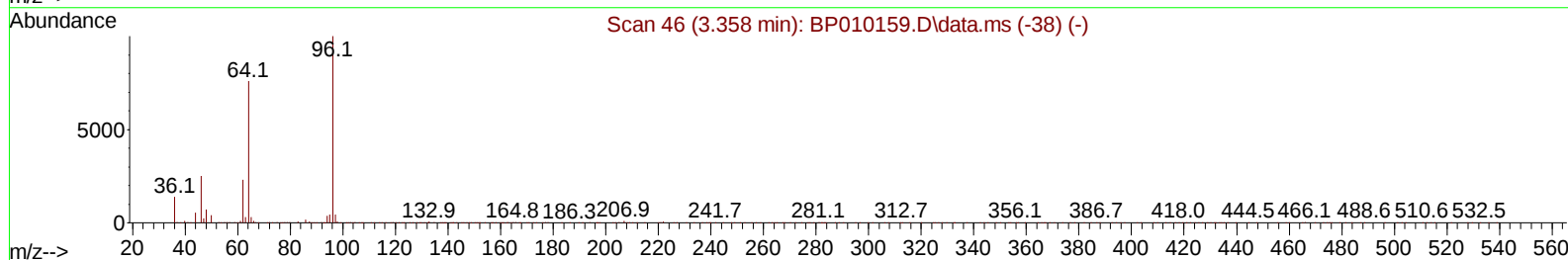
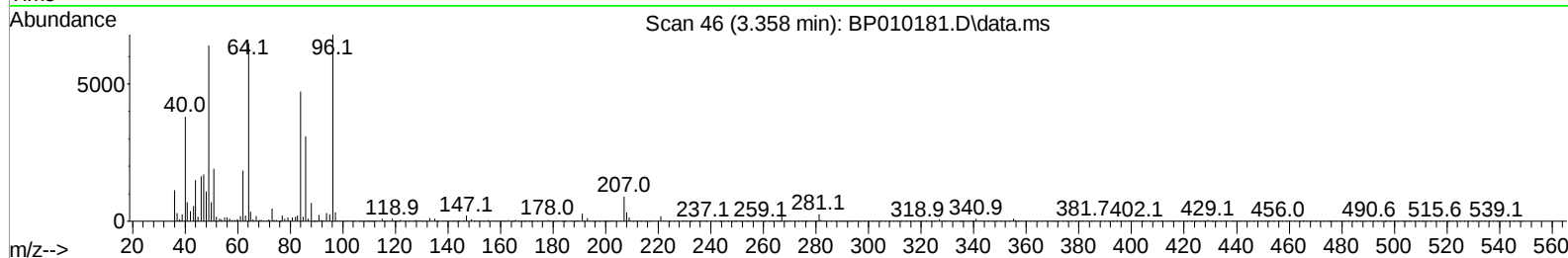
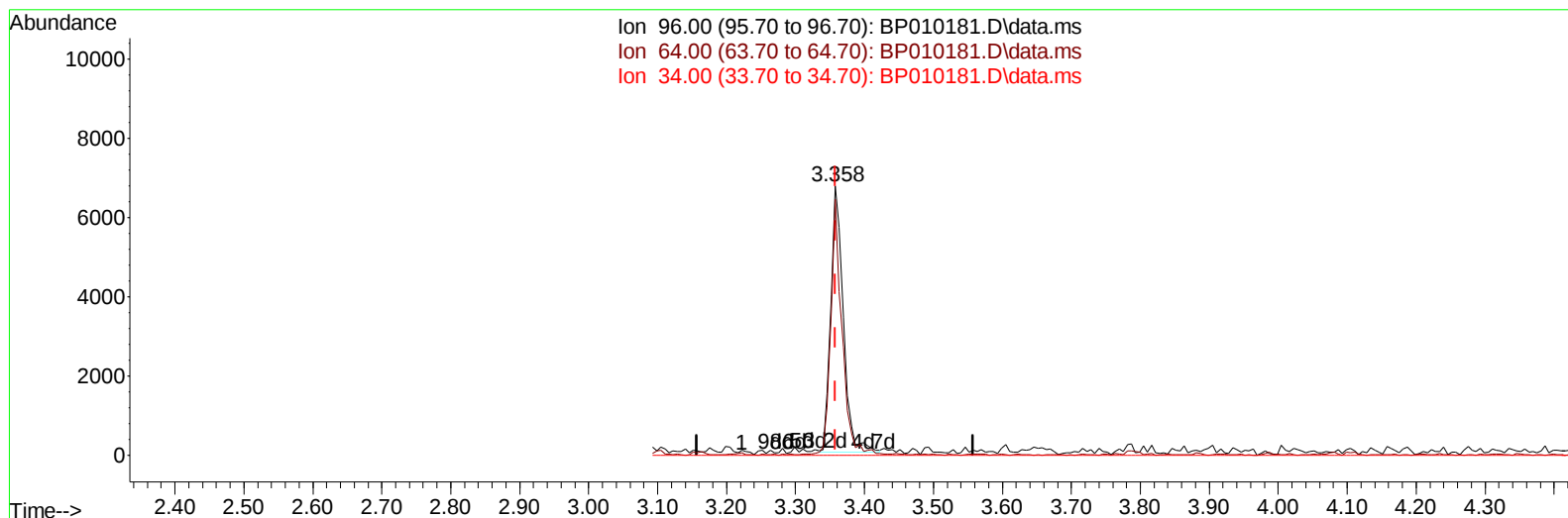
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TIC: BP010181.D\data.ms

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

3.358min (-0.000) 2.76 ng/uL m

response 8592

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	95.10#
34.00	0.00	0.00
0.00	0.00	0.00

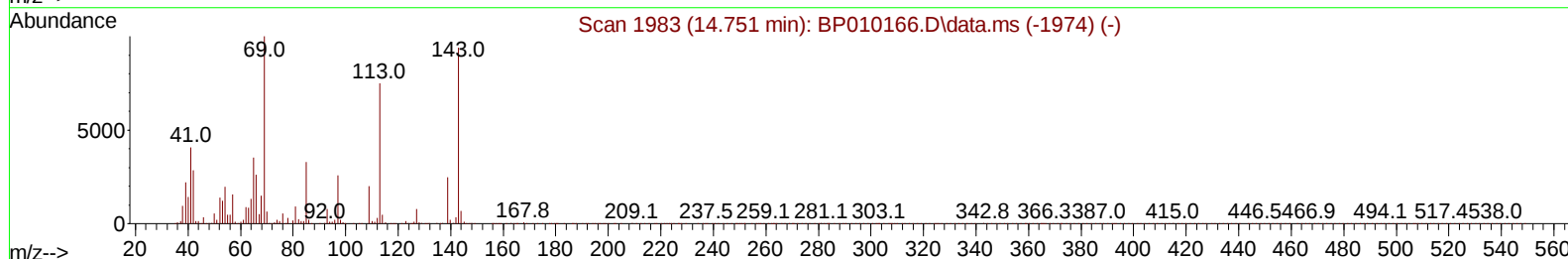
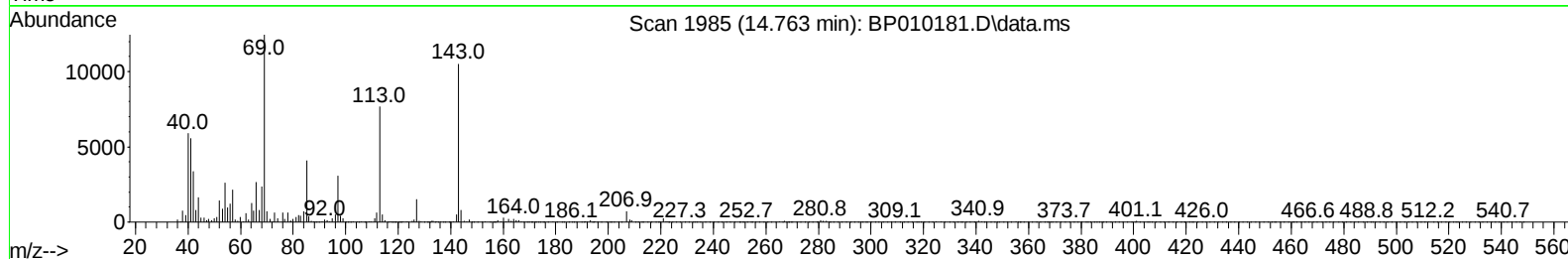
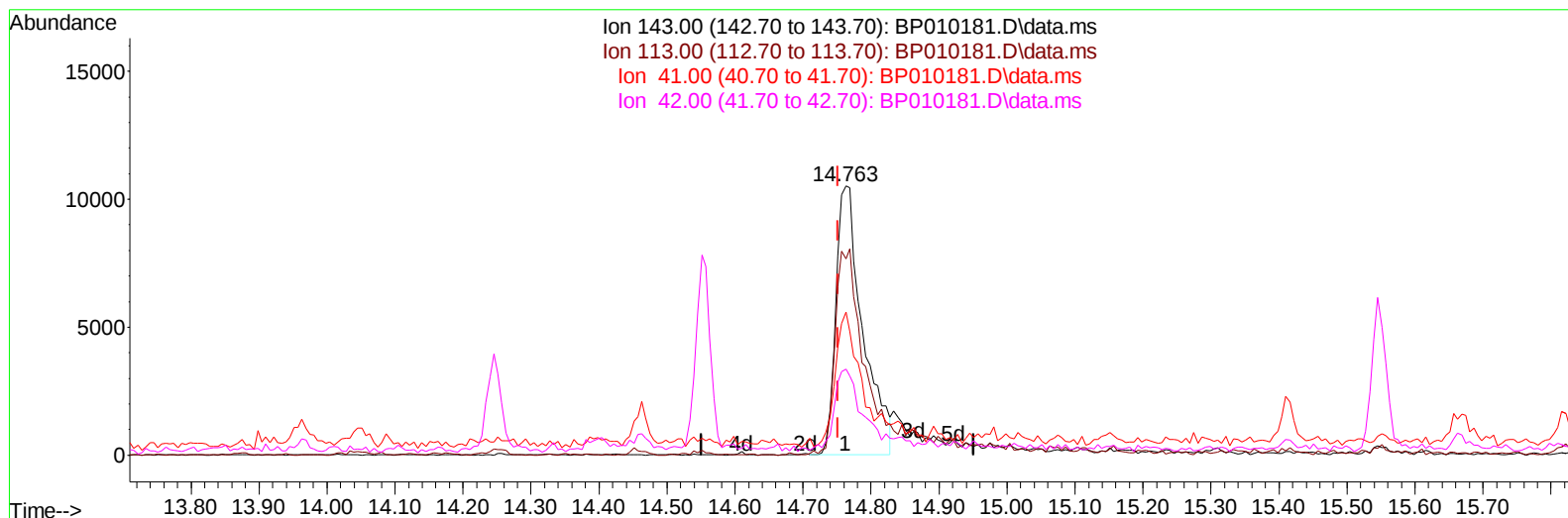
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TIC: BP010181.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 5.98 ng/u1

response 28979

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.05#
41.00	23.20	53.09#
42.00	18.00	32.07#

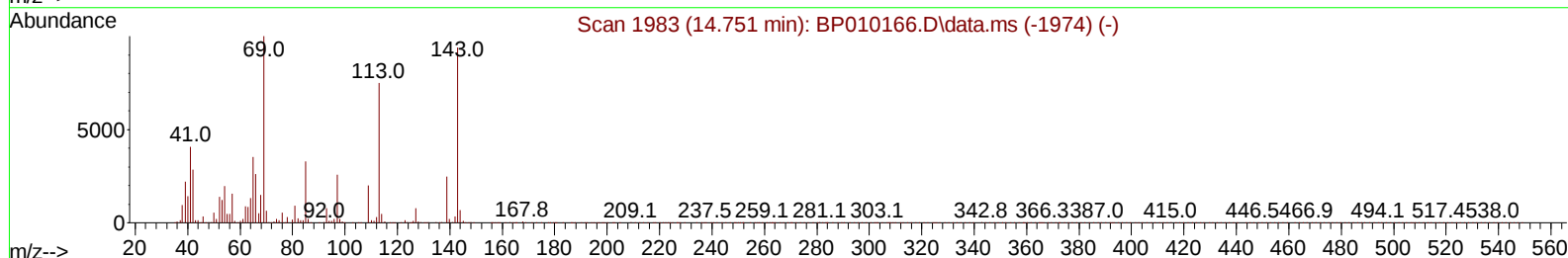
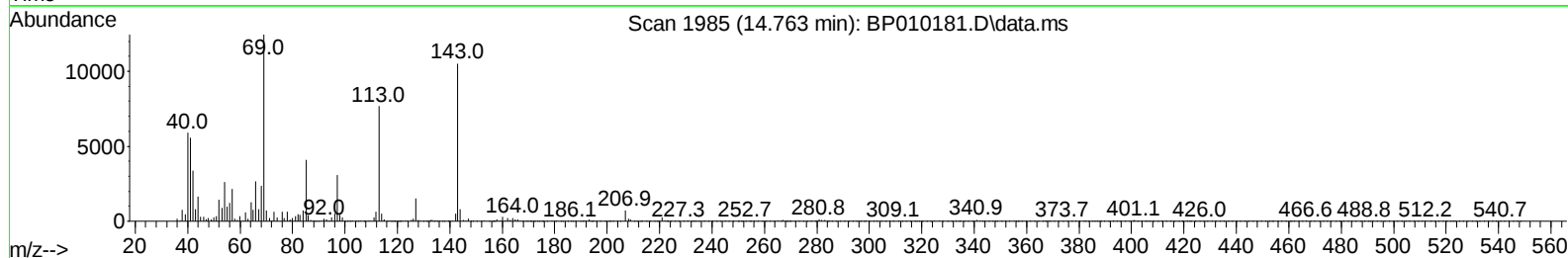
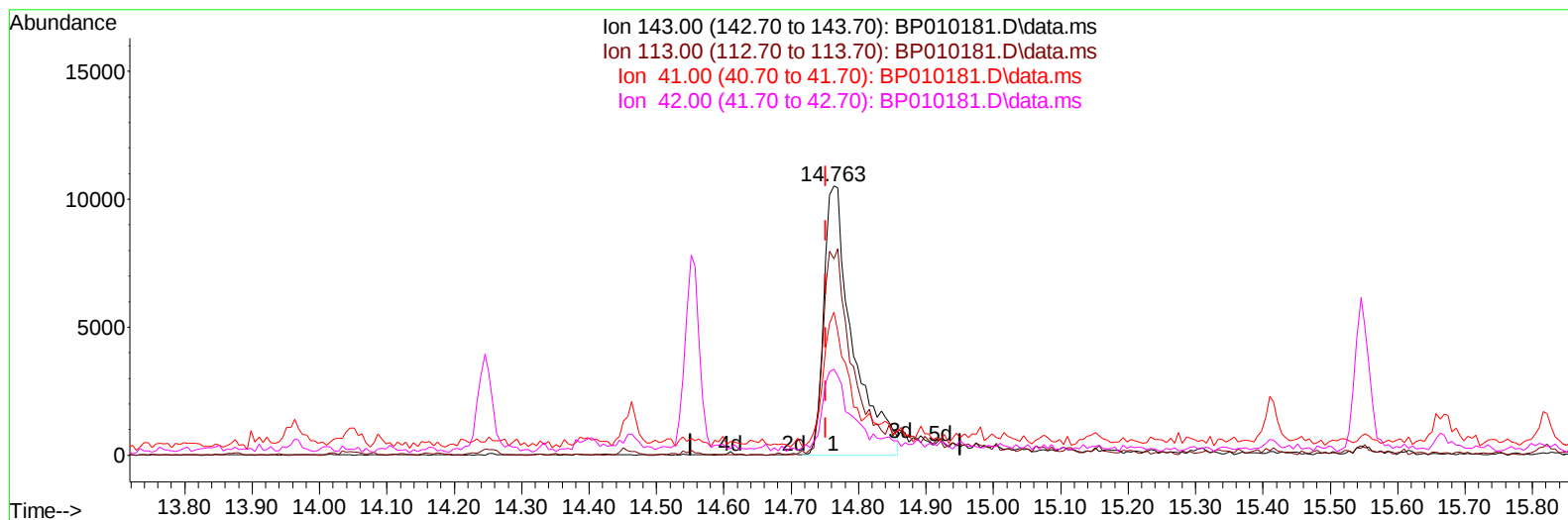
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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP010181.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.012) 6.45 ng/ul m

response 31230

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.05#
41.00	23.20	53.09#
42.00	18.00	32.07#

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.922	152		135784	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136		557188	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164		338213	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188		632555	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240		547273	20.000	ng/ul	# 0.00
88) Perylene-d12	23.833	264		529610	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.358	96		8592m	2.759	ng/uL	0.00
4) Pyridine-d5	3.787	84		37813	4.317	ng/ul	0.01
7) Phenol-d5	7.087	99		155112	12.559	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67		110836	14.956	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132		132619	14.499	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113		127292	12.439	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128		66311	14.963	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143		72077	14.892	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165		116035	12.315	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131		157690	11.672	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166		383380	14.572	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160		445231	13.983	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143		31230m	6.448	ng/ul	0.01
60) Fluorene-d10	15.545	176		328109	14.498	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200		25138	6.285	ng/ul	0.00
73) Anthracene-d10	17.404	188		469739	15.350	ng/ul	0.00
81) Pyrene-d10	19.633	212		569348	18.293	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264		471199	16.971	ng/ul	0.00

Target Compounds Qvalue

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

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