

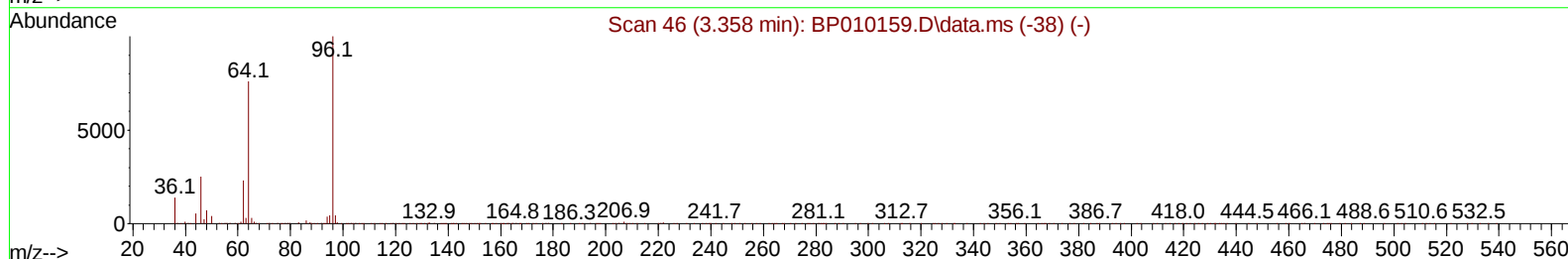
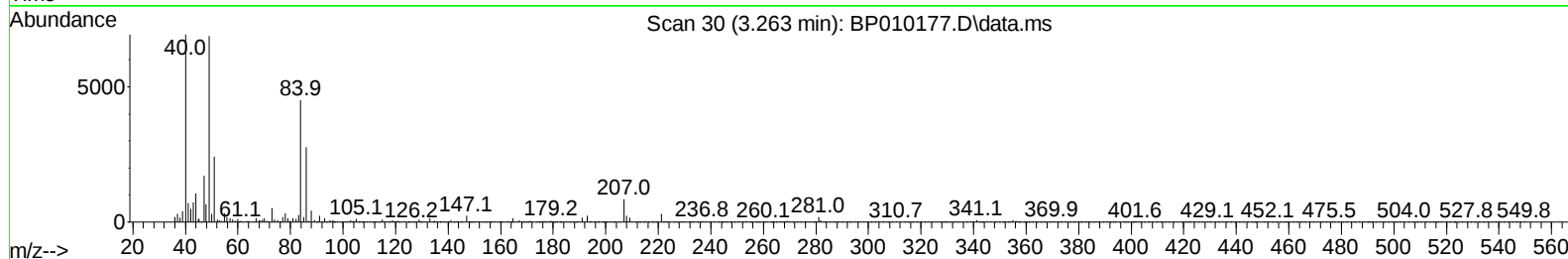
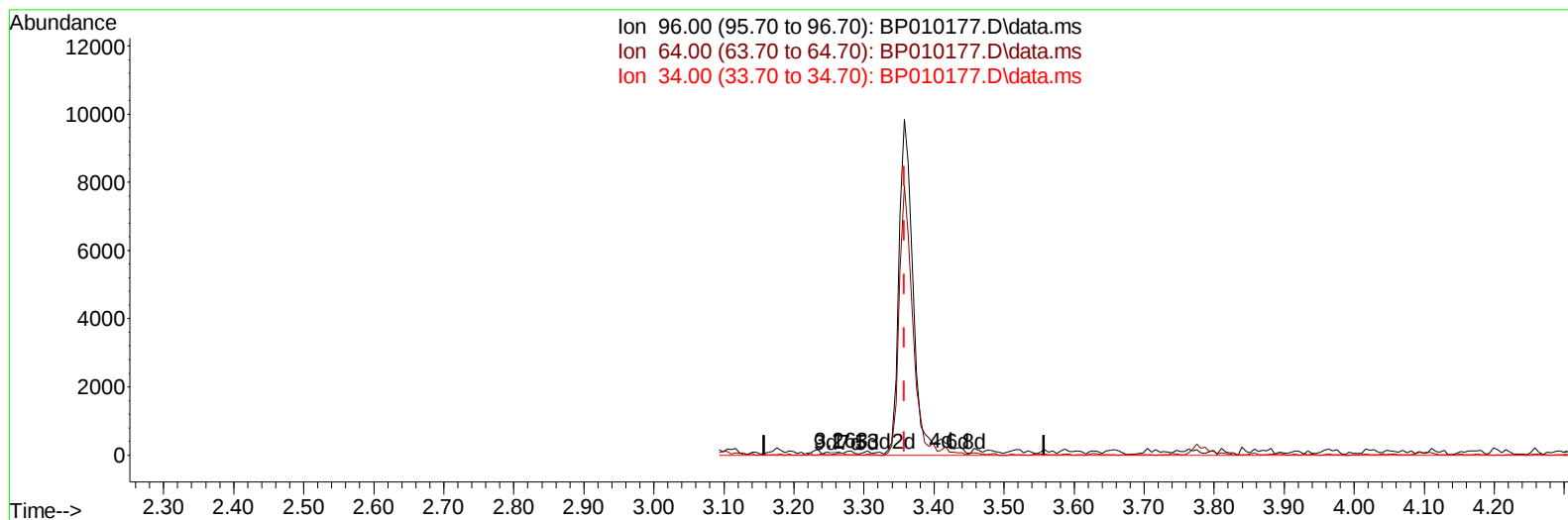
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010177.D
 Acq On : 02 May 2022 17:32
 Operator : CG/JU
 Sample : N2562-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW478

Manual IntegrationsAPPROVED

Quant Time: May 03 01:37:14 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: BP010177.D\data.ms

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

3.263min (-0.094) 0.01 ng/uL

response 18

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	62.20
34.00	0.00	0.00
0.00	0.00	0.00

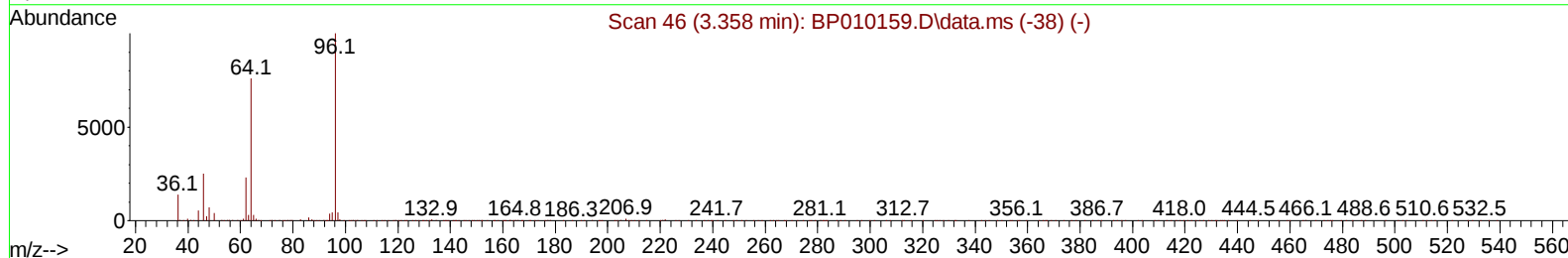
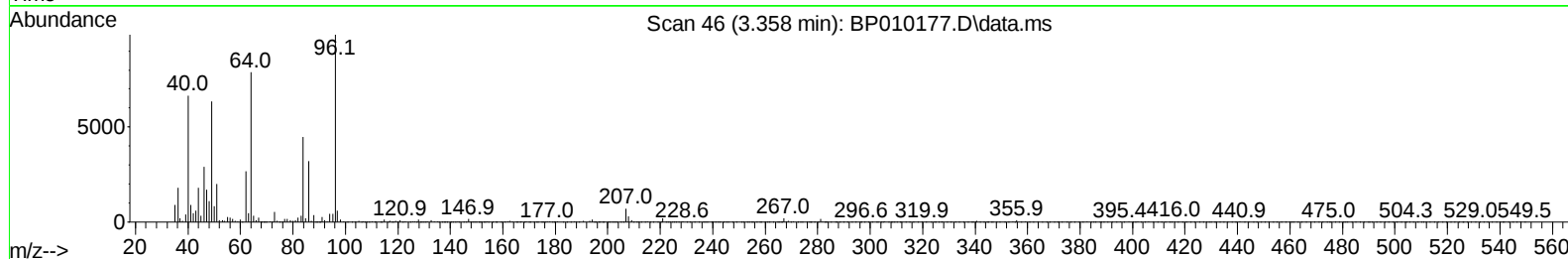
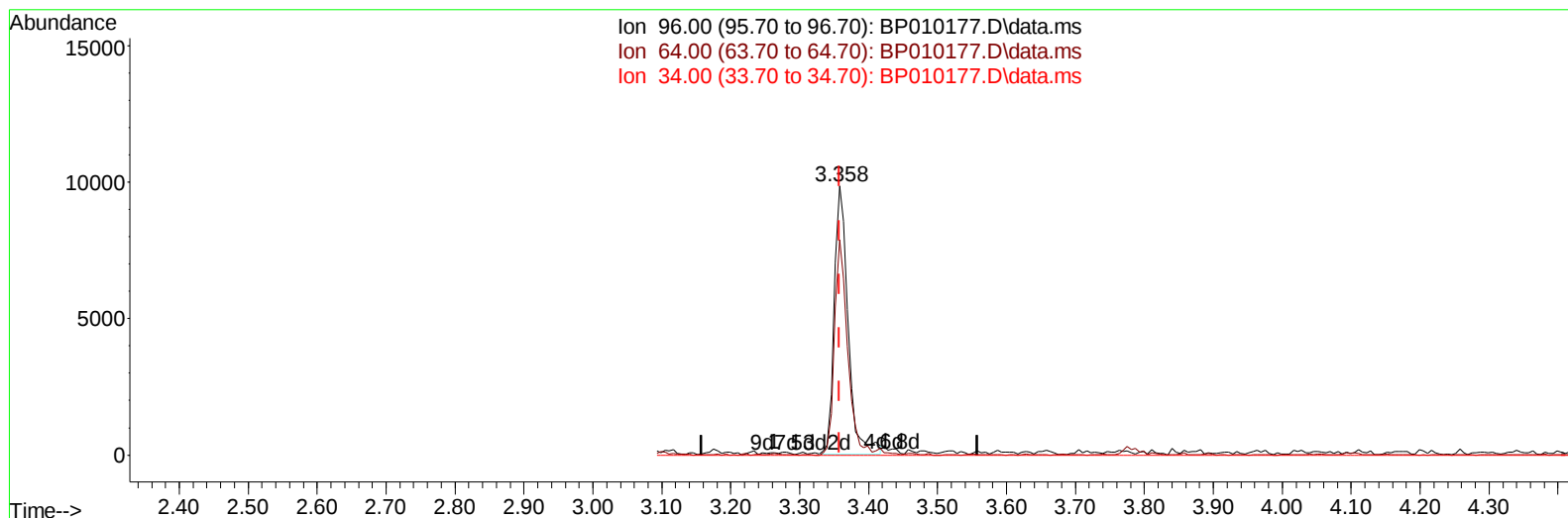
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(3) 1,4-Dioxane-d8 (S)

3.358min (-0.000) 4.20 ng/uL m

response 13809

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.04#
34.00	0.00	0.00
0.00	0.00	0.00

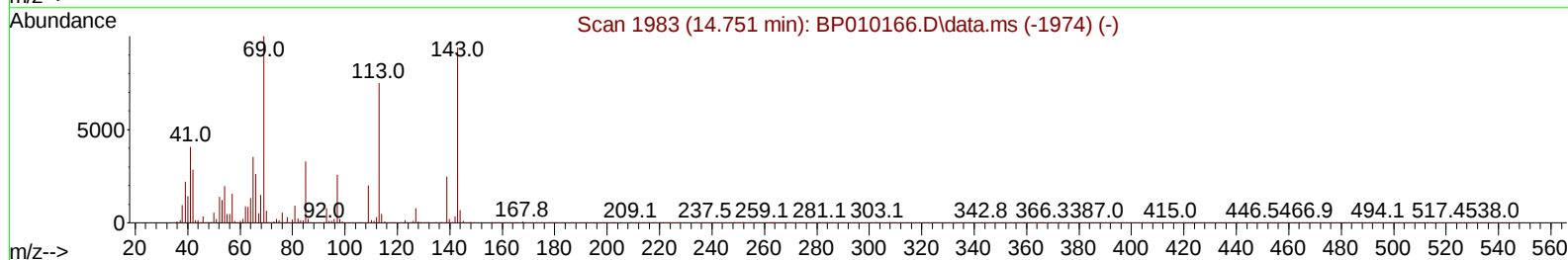
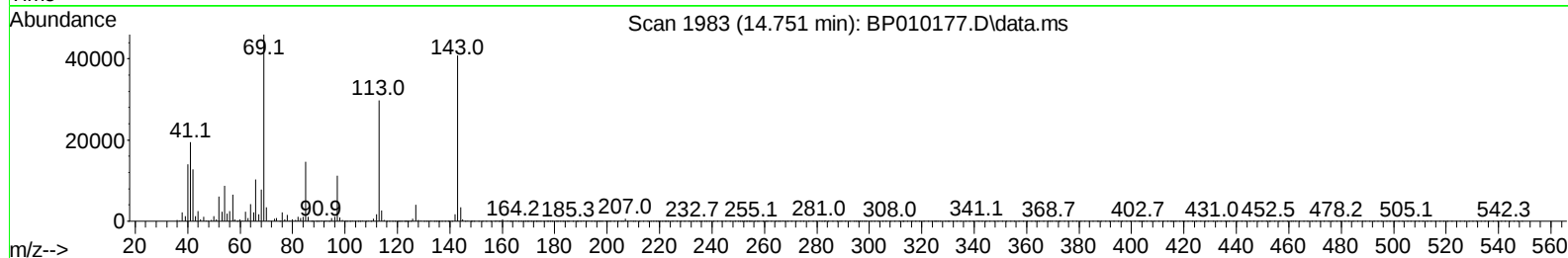
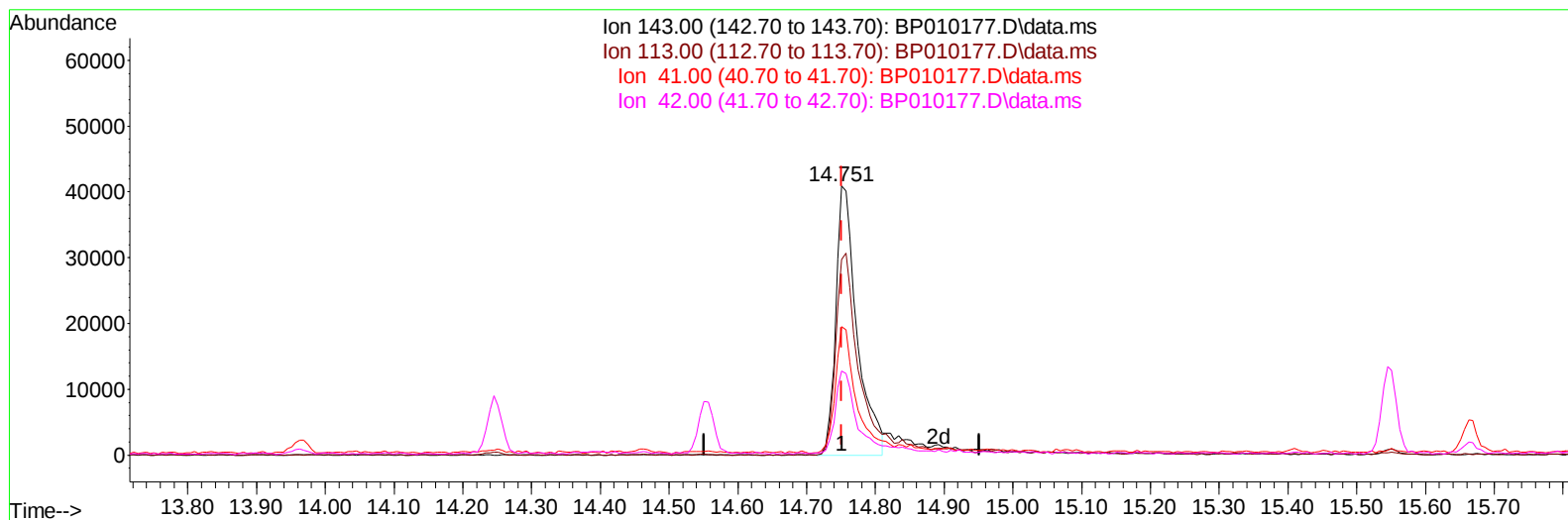
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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TIC: BP010177.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 15.66 ng/u1

response 87693

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.86#
41.00	23.20	47.69#
42.00	18.00	31.26#

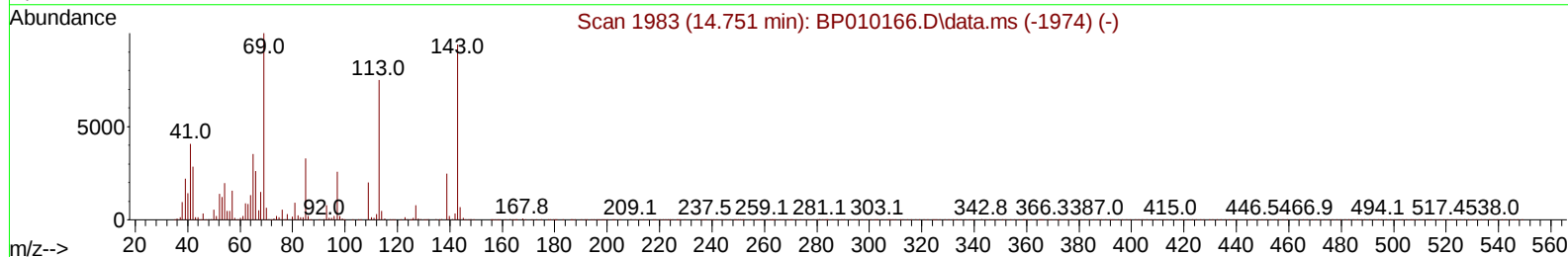
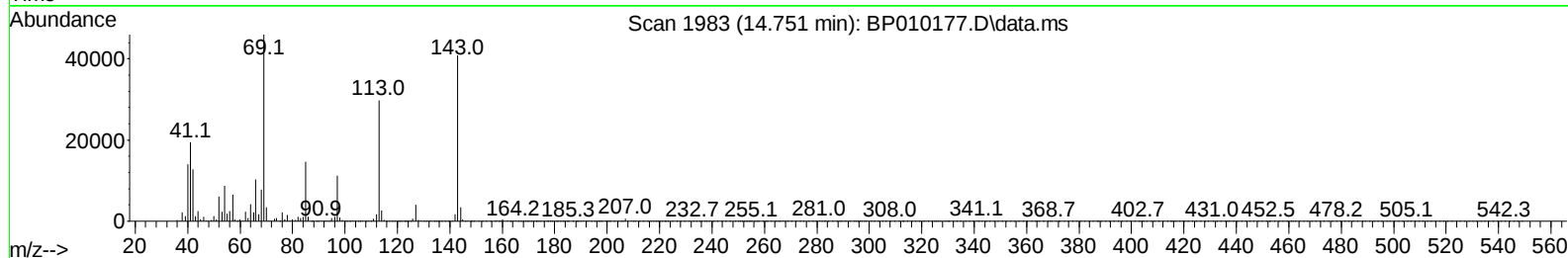
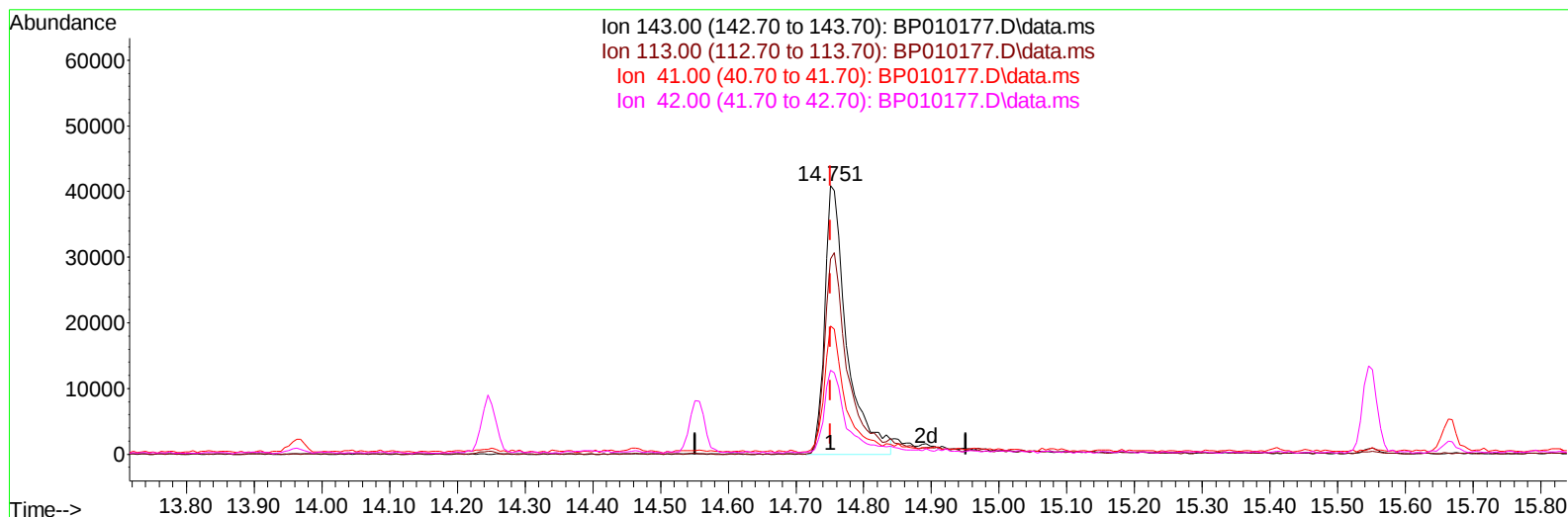
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW478

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Supervised By :Yogesh Patel 05/05/2022

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TIC: BP010177.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 16.56 ng/ul m

response 92758

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.86#
41.00	23.20	47.69#
42.00	18.00	31.26#

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.922	152		143379	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136		613701	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164		391146	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188		726944	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.392	240		553631	20.000	ng/ul	0.00
88) Perylene-d12	23.833	264		525770	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.358	96		13809m	4.200	ng/uL	0.00
4) Pyridine-d5	3.775	84		172003	18.595	ng/ul	0.00
7) Phenol-d5	7.087	99		286774	21.989	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67		201658	25.769	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132		235612	24.394	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113		257624	23.842	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128		130428	26.721	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143		142555	26.742	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165		245862	23.691	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131		360449	24.224	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166		908425	29.856	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160		1024621	27.824	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143		92758m	16.560	ng/ul	0.00
60) Fluorene-d10	15.545	176		749224	28.626	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200		81098	17.643	ng/ul	0.00
73) Anthracene-d10	17.404	188		1053141	29.945	ng/ul	0.00
81) Pyrene-d10	19.639	212		1110238	35.262	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264		858746	31.155	ng/ul	0.00

Target Compounds Qvalue

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

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