

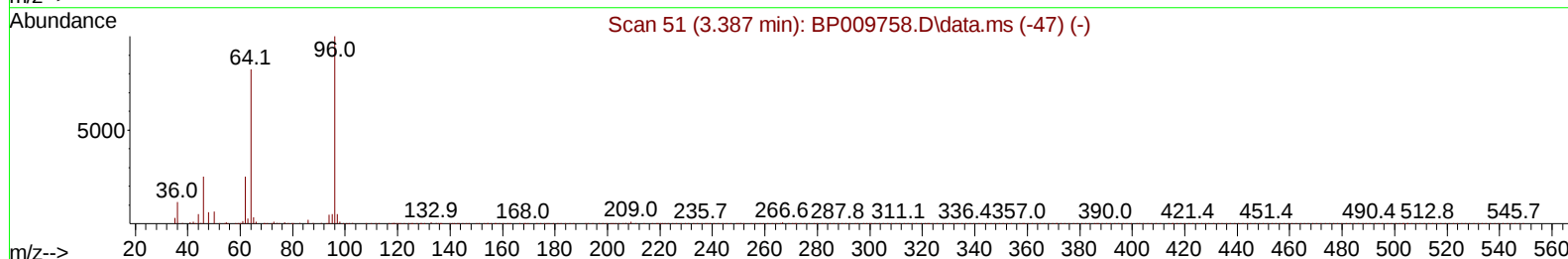
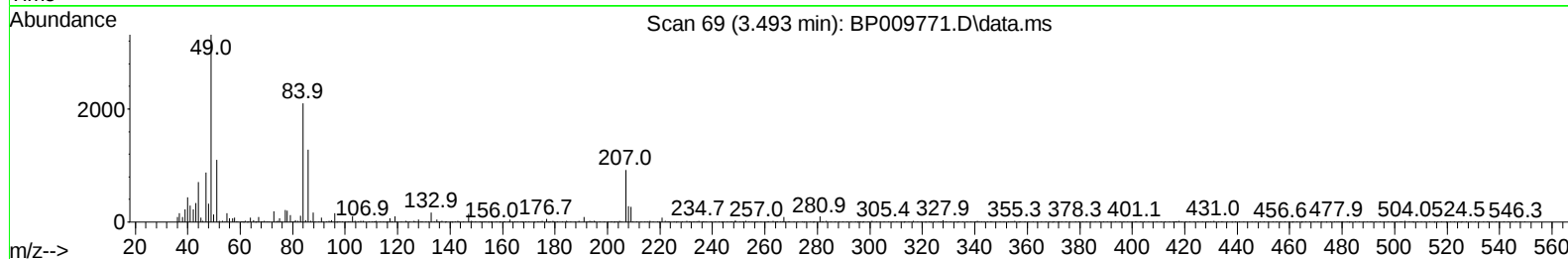
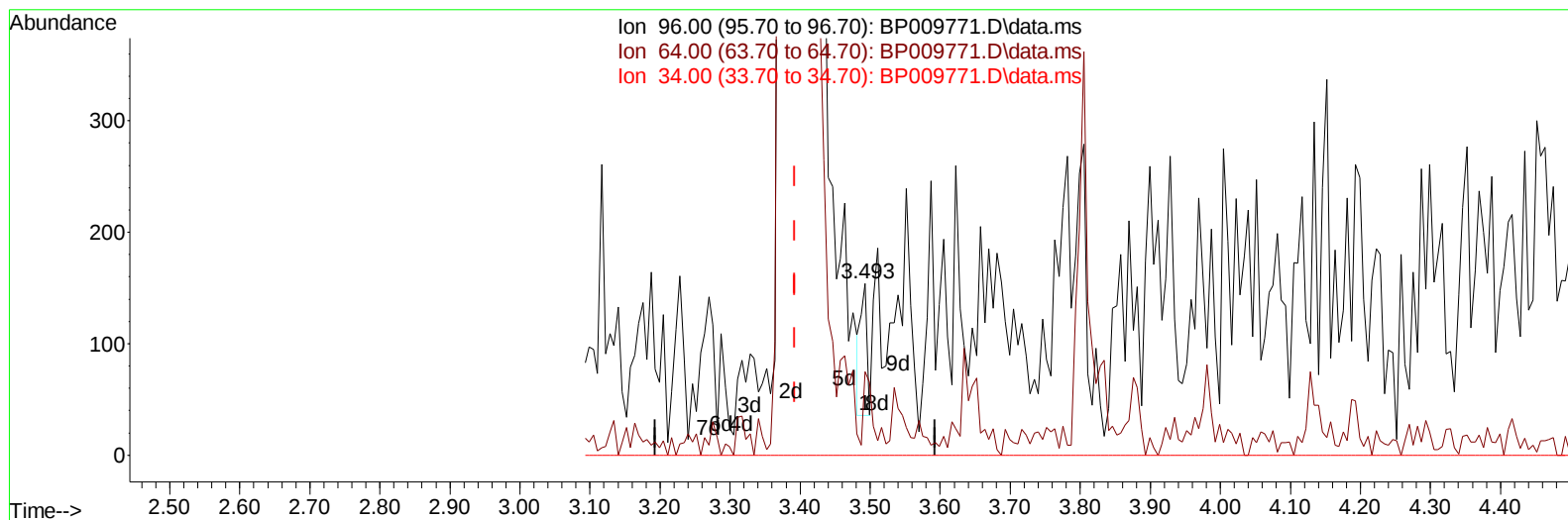
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
 Data File : BP009771.D
 Acq On : 12 Apr 2022 02:38
 Operator : CG/JU
 Sample : N2338-19
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J83

Manual IntegrationsAPPROVED

Quant Time: Apr 12 03:49:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 07 19:16:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/13/2022
 Supervised By :Yogesh Patel 04/13/2022



TIC: BP009771.D\data.ms

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

3.493min (+ 0.100) 0.02 ng/uL

response 73

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.39
34.00	0.00	0.00
0.00	0.00	0.00

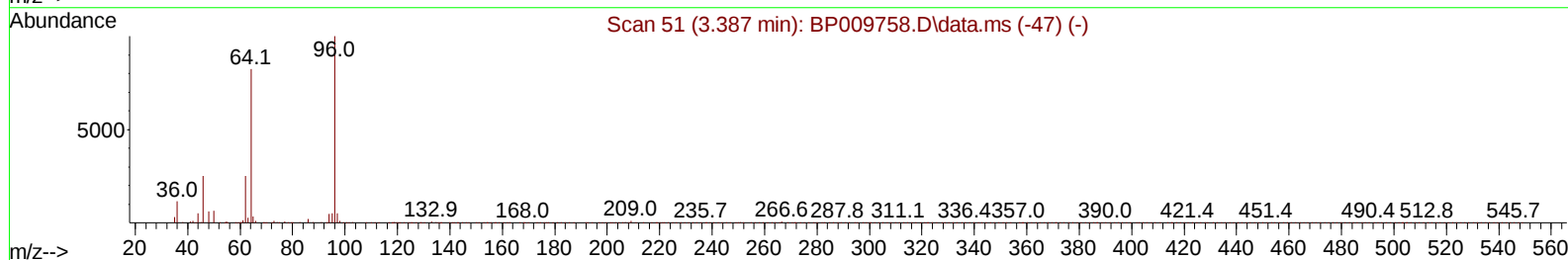
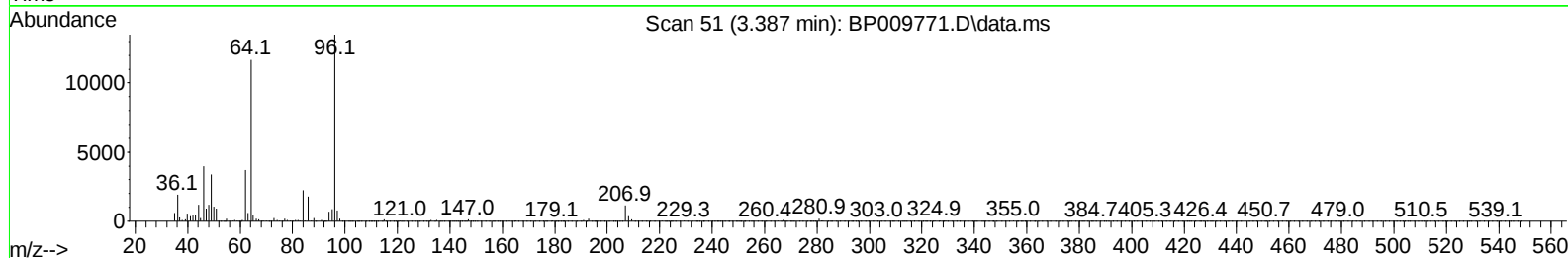
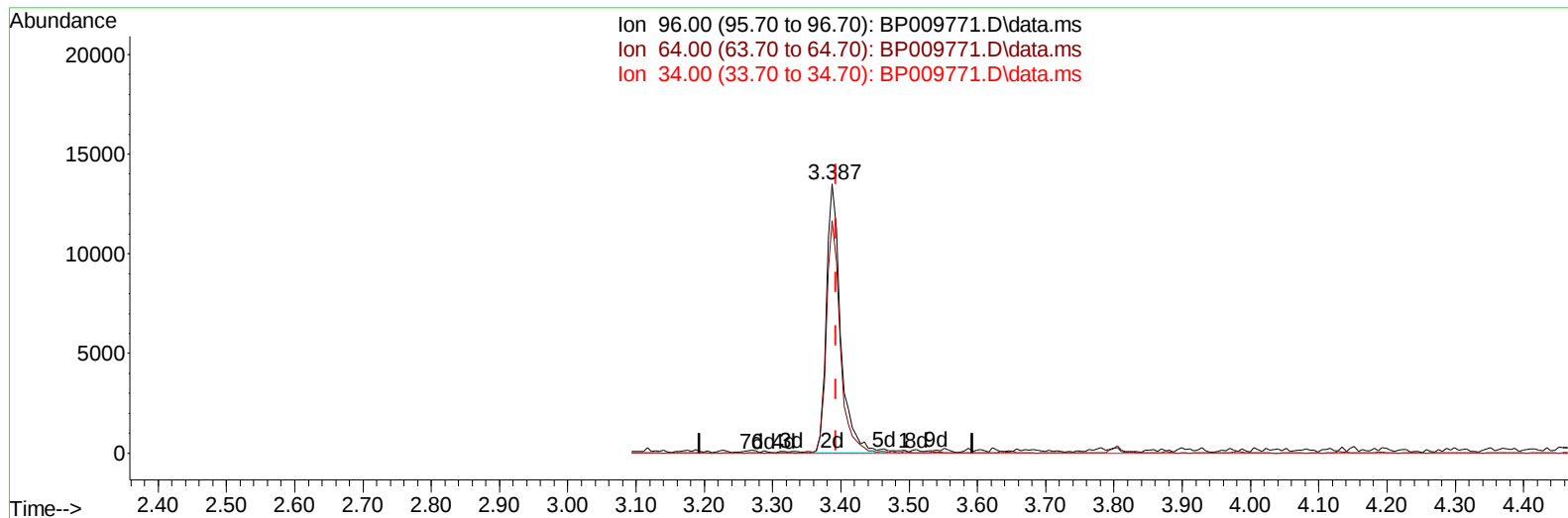
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(3) 1,4-Dioxane-d8 (S)

3.387min (-0.006) 6.52 ng/uL m

response 19407

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.46#
34.00	0.00	0.00
0.00	0.00	0.00

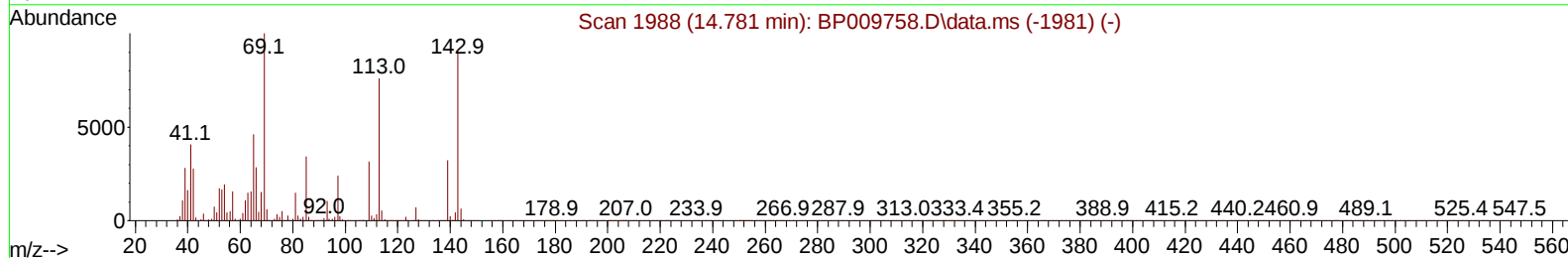
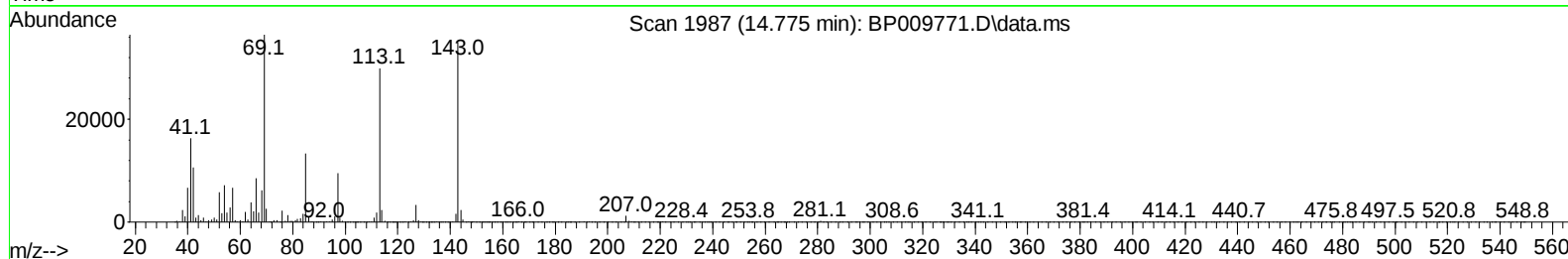
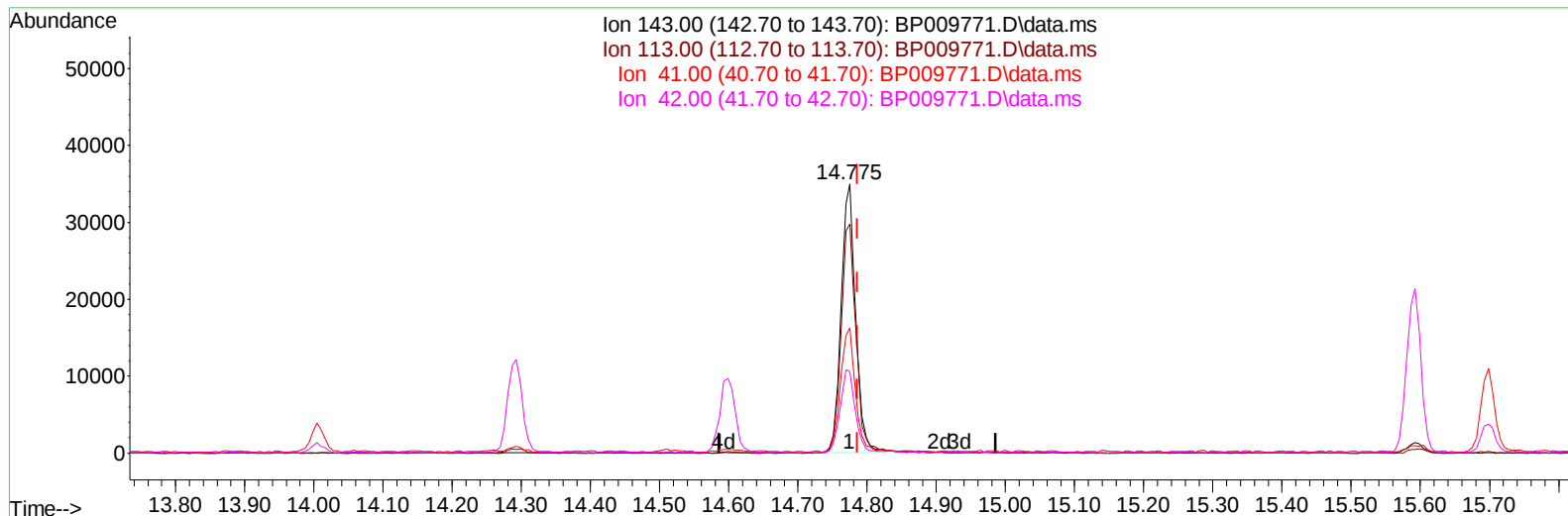
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TIC: BP009771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 8.06 ng/u1

response 50726

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.32#
41.00	23.20	46.51#
42.00	18.00	30.21#

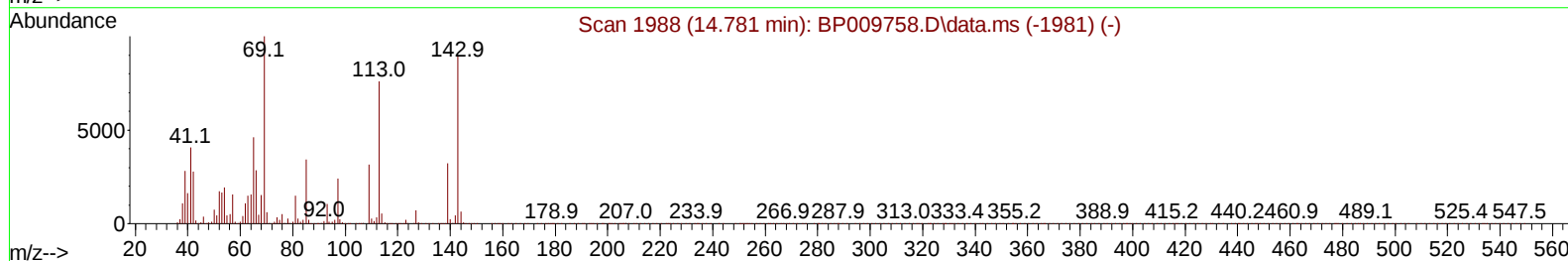
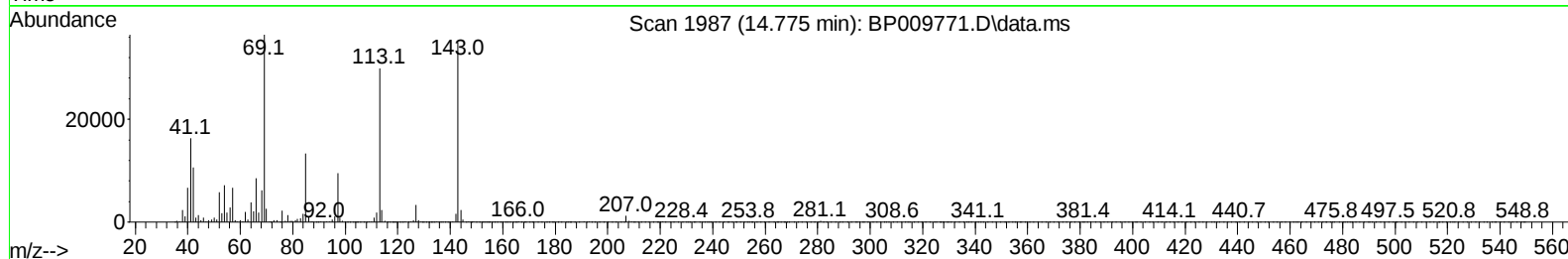
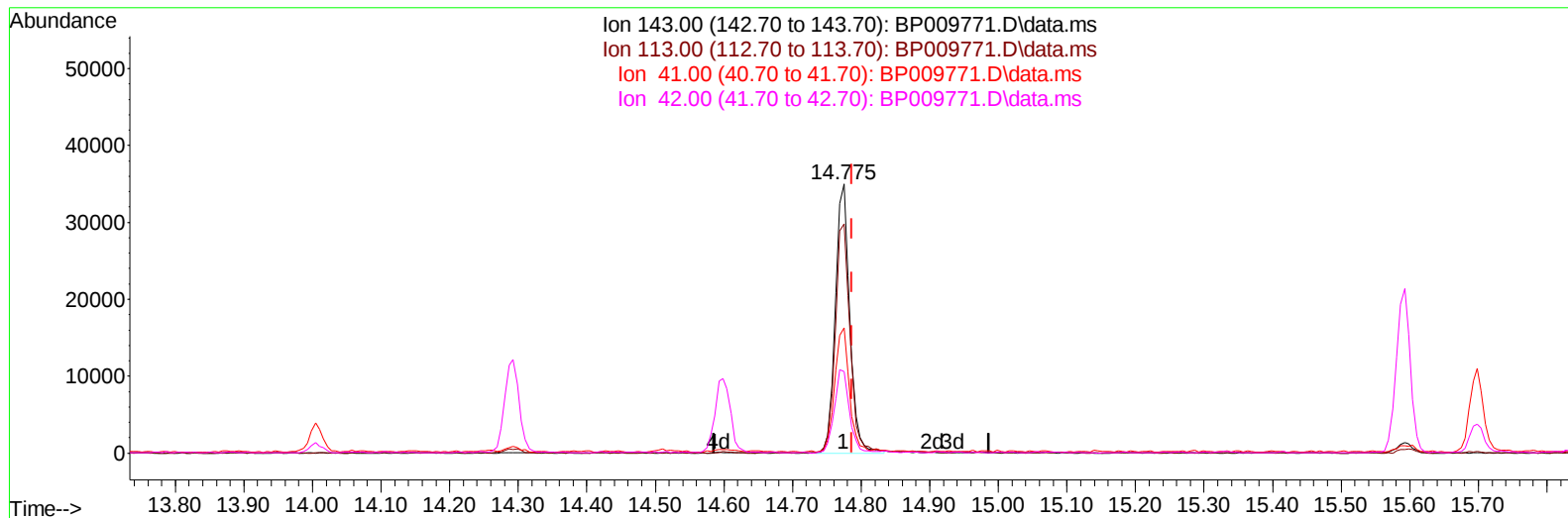
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Instrument :
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TIC: BP009771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.775min (-0.012) 8.13 ng/ul m

response 51170

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.32#
41.00	23.20	46.51#
42.00	18.00	30.21#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	132761	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.775	136	602619	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.598	164	441653	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.351	188	973892	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.427	240	987389	20.000	ng/ul	#-0.02
88) Perylene-d12	23.898	264	922448	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	19407m	6.524	ng/uL	0.00
4) Pyridine-d5	3.805	84	116552	14.211	ng/ul	0.00
7) Phenol-d5	7.122	99	94558	8.113	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	243949	35.156	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.499	132	218844	24.891	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	175990	18.087	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	158025	36.364	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.852	143	153782	32.278	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	267942	26.876	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.905	131	439852	30.976	ng/ul	-0.01
46) Dimethylphthalate-d6	14.004	166	1226906	35.365	ng/ul	-0.02
49) Acenaphthylene-d8	14.293	160	1342791	32.671	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.775	143	51170m	8.133	ng/ul	-0.01
60) Fluorene-d10	15.592	176	1068610	36.616	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.698	200	168204	29.155	ng/ul	-0.01
73) Anthracene-d10	17.451	188	1843685	39.539	ng/ul	-0.01
81) Pyrene-d10	19.675	212	2225005	41.260	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.739	264	1961717	40.837	ng/ul	-0.02

Target Compounds Qvalue

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

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