

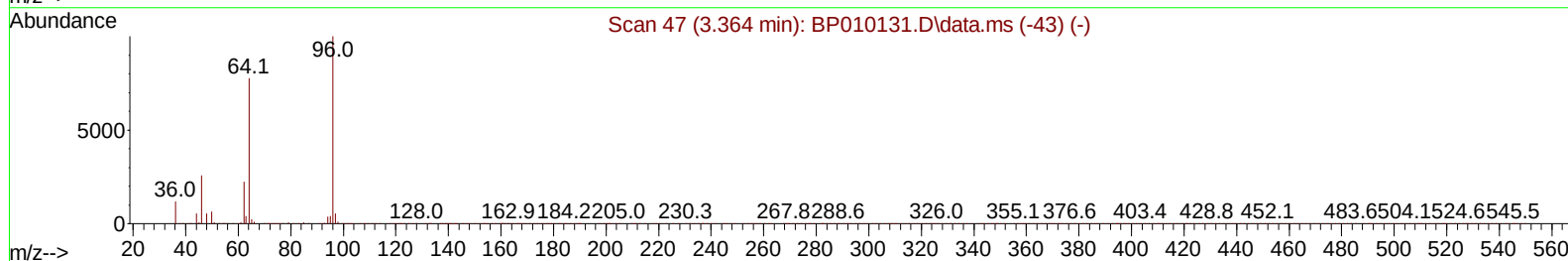
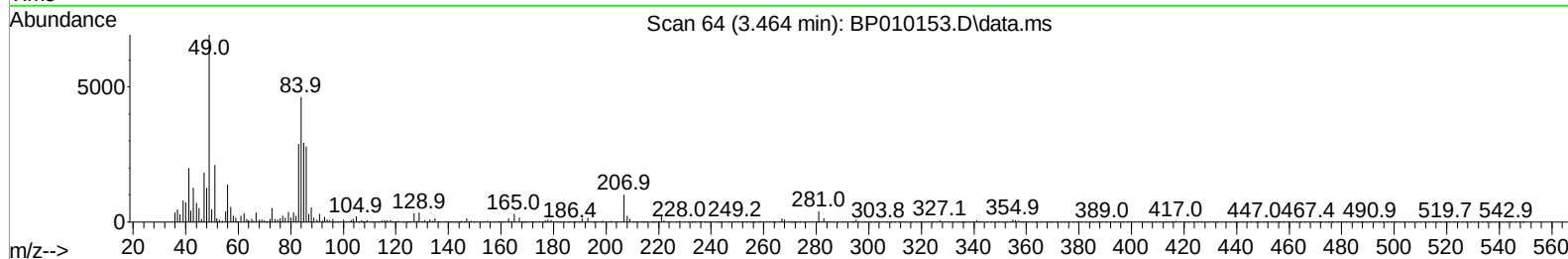
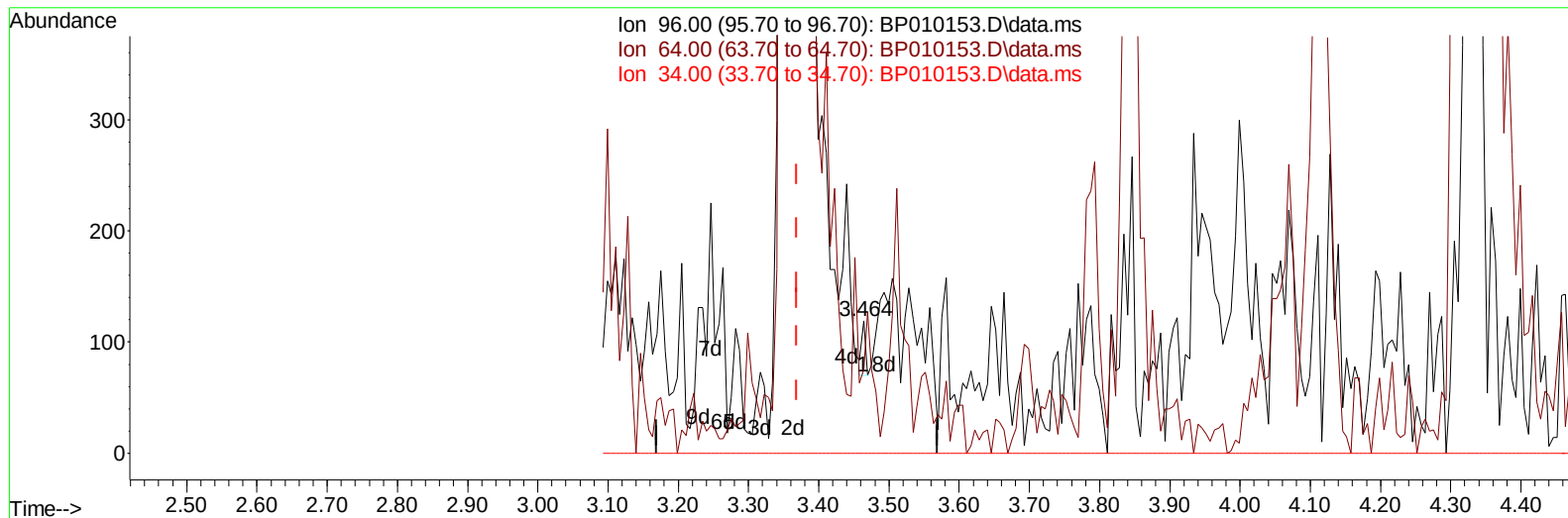
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
 Data File : BP010153.D
 Acq On : 29 Apr 2022 16:09
 Operator : CG/JU
 Sample : N2587-18
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXET2

Manual IntegrationsAPPROVED

Quant Time: Apr 30 00:30:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/02/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010153.D\data.ms

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

3.464min (+ 0.095) 0.01 ng/uL

response 17

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	57.69
34.00	0.00	0.00
0.00	0.00	0.00

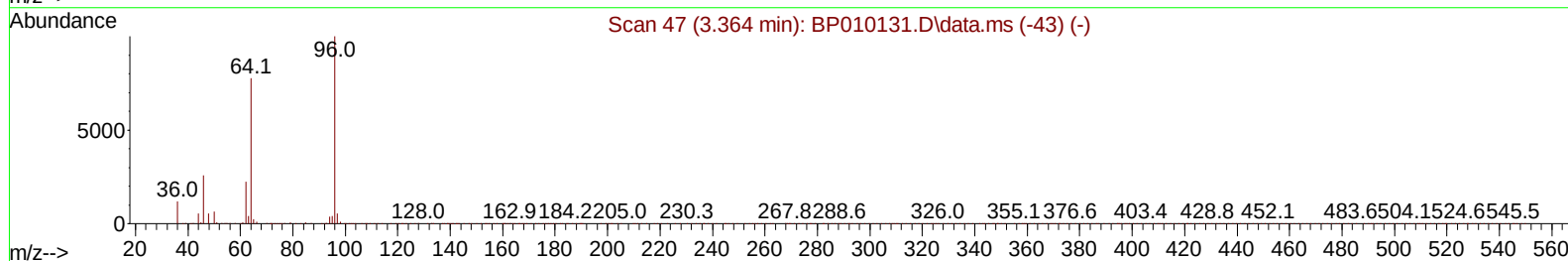
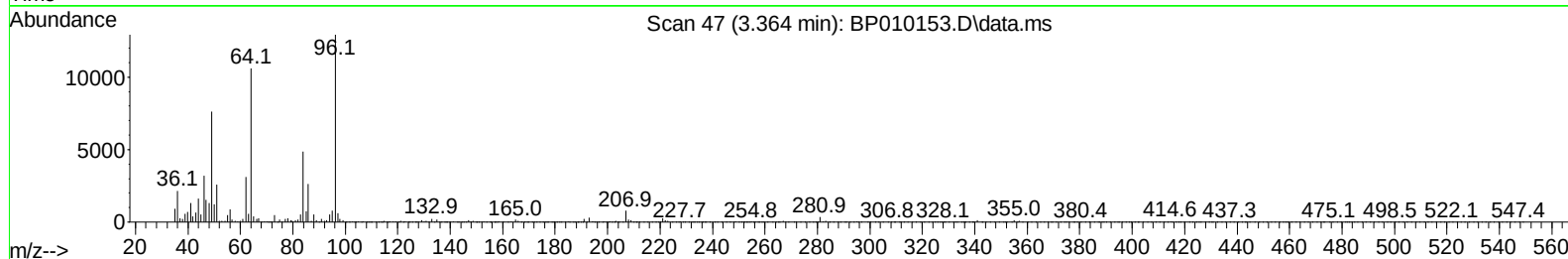
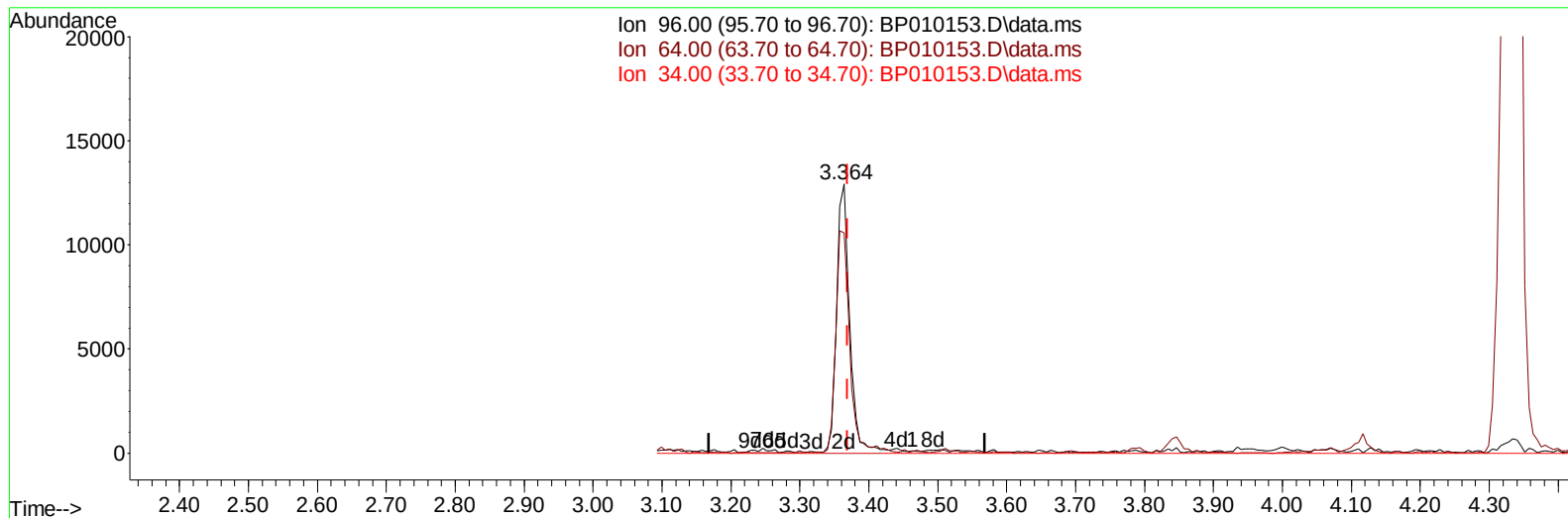
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(3) 1,4-Dioxane-d8 (S)

3.364min (-0.005) 5.50 ng/uL m

response 16864

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.89#
34.00	0.00	0.00
0.00	0.00	0.00

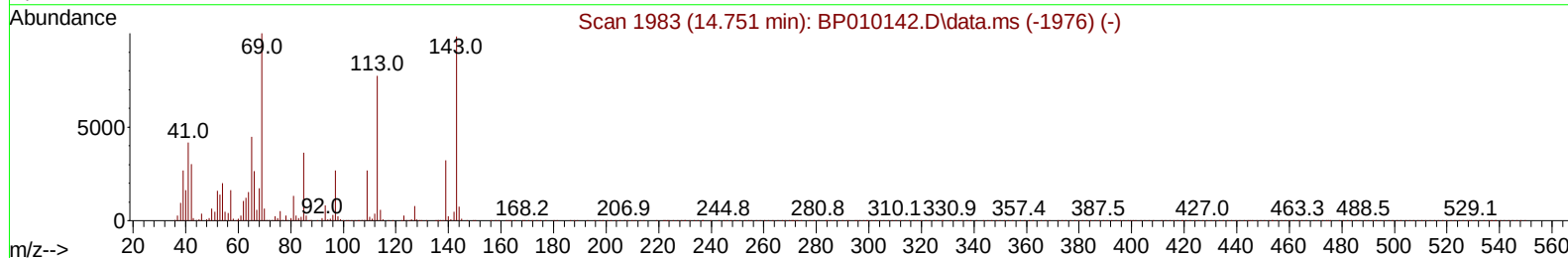
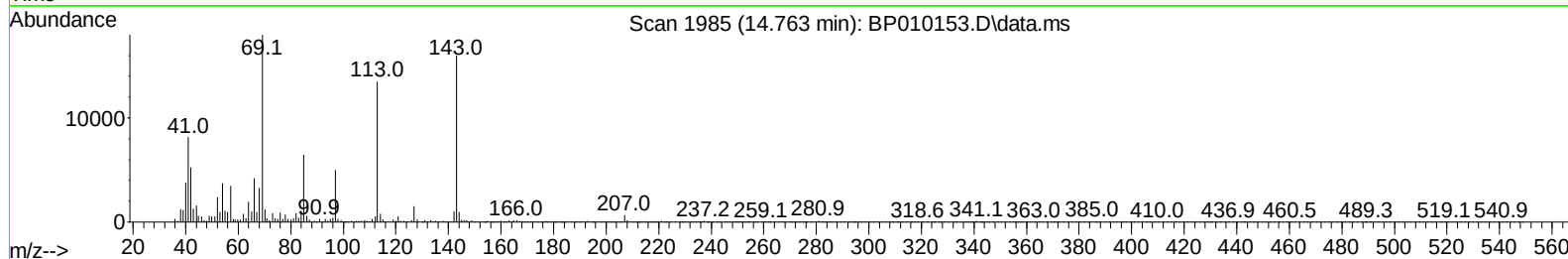
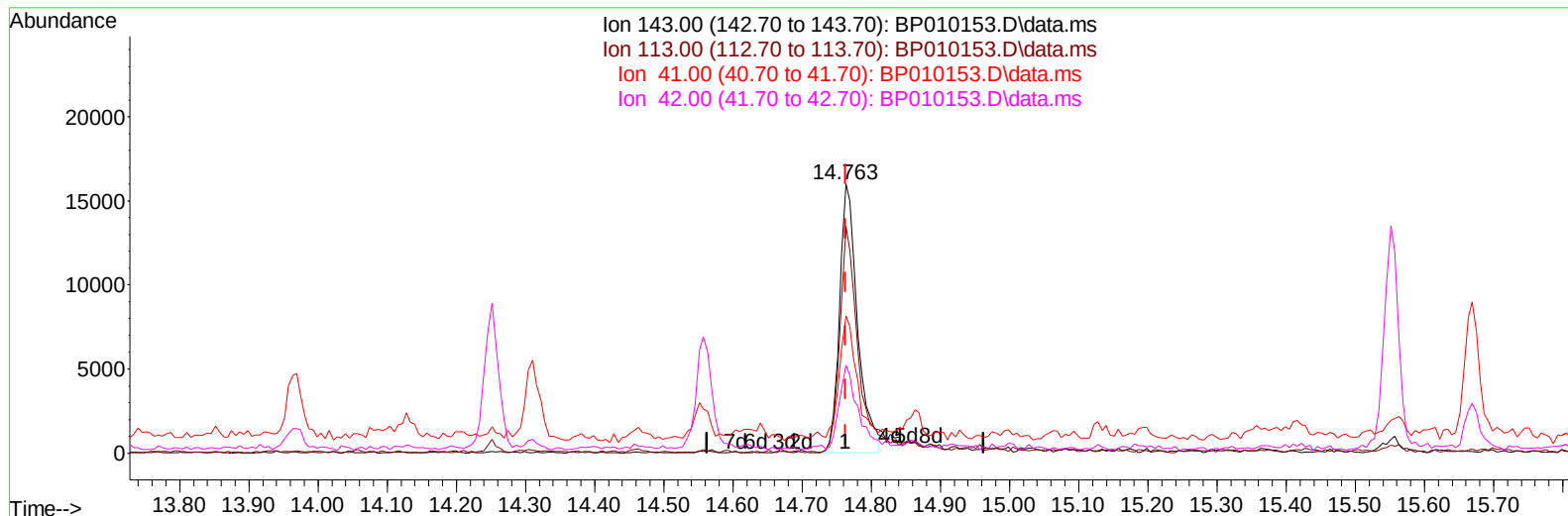
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TIC: BP010153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 6.43 ng/ul

response 29139

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.33#
41.00	23.20	51.00#
42.00	18.00	32.70#

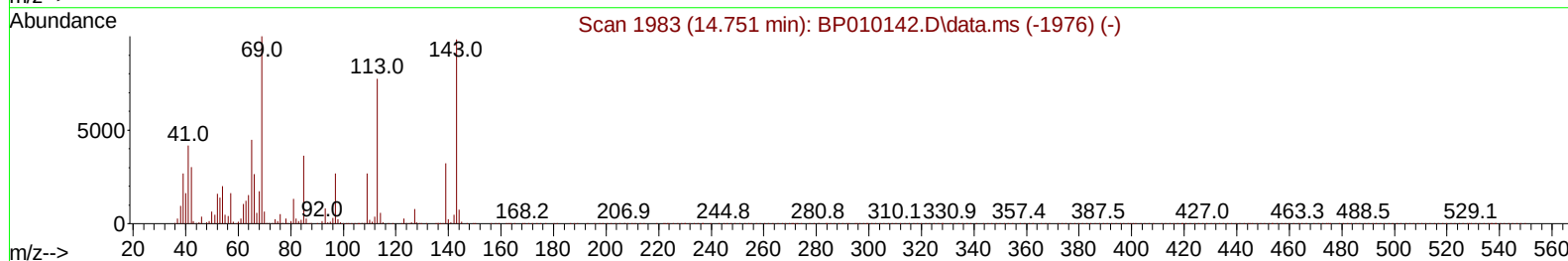
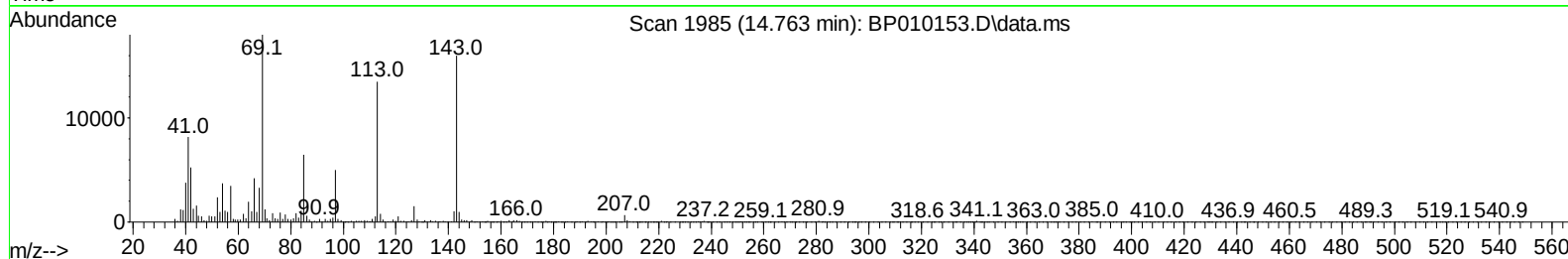
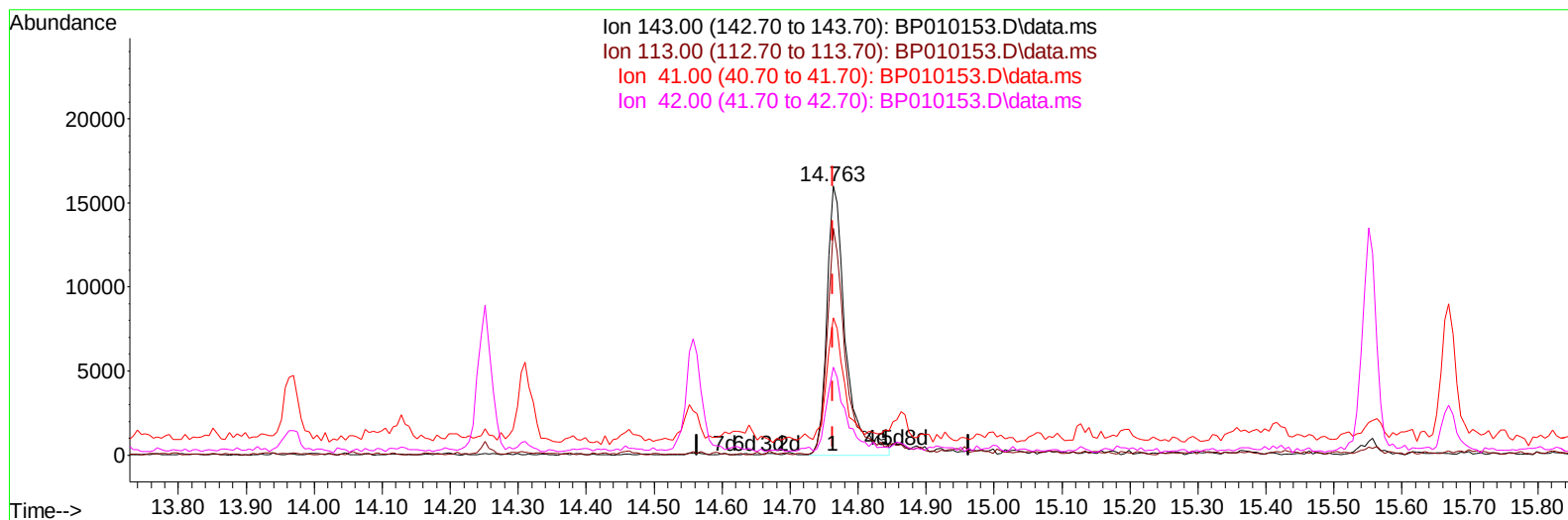
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Instrument :
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TIC: BP010153.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 6.86 ng/ul m

response 31106

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.33#
41.00	23.20	51.00#
42.00	18.00	32.70#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	133652	20.000	ng/ul	#-0.01
20) Naphthalene-d8	10.728	136	515884	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.557	164	316643	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.310	188	625994	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.392	240	591501	20.000	ng/ul	#-0.02
88) Perylene-d12	23.845	264	595884	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	16864m	5.502	ng/uL	0.00
4) Pyridine-d5	3.787	84	96354	11.175	ng/ul	0.00
7) Phenol-d5	7.093	99	104947	8.633	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	235146	32.235	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.458	132	240320	26.693	ng/ul	-0.01
15) 4-Methylphenol-d8	8.634	113	188777	18.742	ng/ul	-0.02
21) Nitrobenzene-d5	9.087	128	140878	34.335	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.810	143	153470	34.248	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	252782	28.977	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	291496	23.304	ng/ul	-0.02
46) Dimethylphthalate-d6	13.969	166	862563	35.019	ng/ul	-0.01
49) Acenaphthylene-d8	14.251	160	977737	32.798	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	31106m	6.860	ng/ul	0.00
60) Fluorene-d10	15.551	176	729090	34.411	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200	129162	32.631	ng/ul	-0.01
73) Anthracene-d10	17.410	188	1147185	37.880	ng/ul	-0.01
81) Pyrene-d10	19.645	212	1342330	39.904	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1261870	40.393	ng/ul	-0.01

Target Compounds Qvalue

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

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