

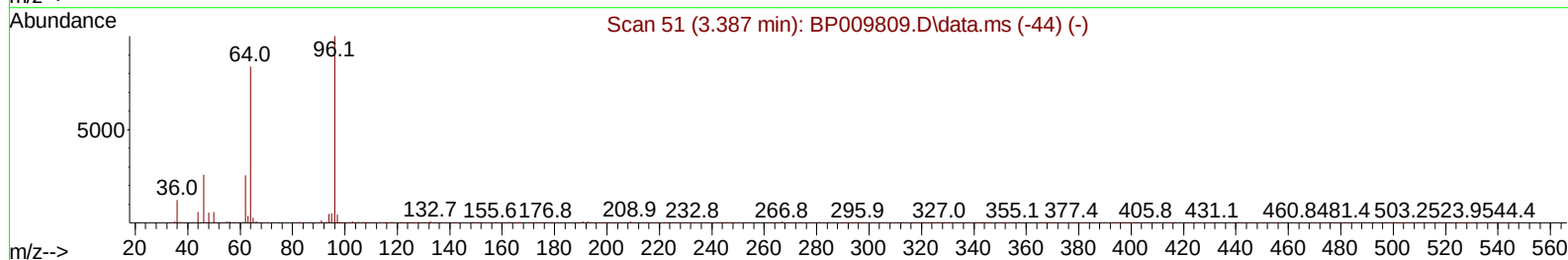
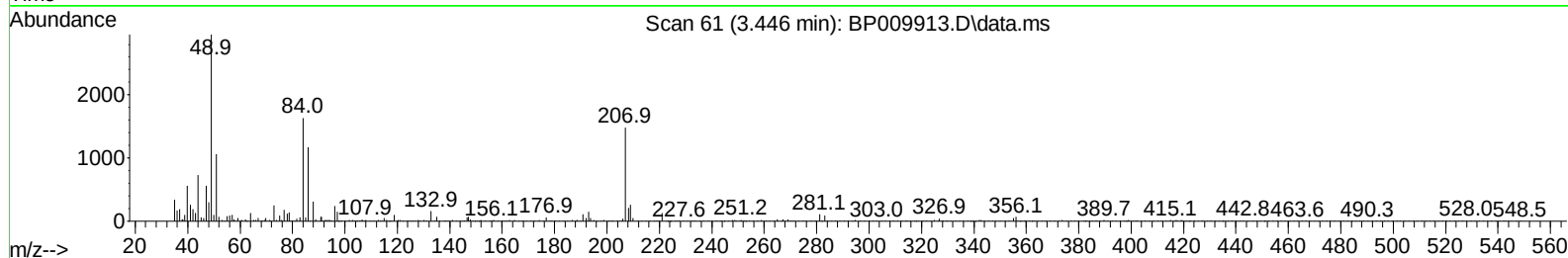
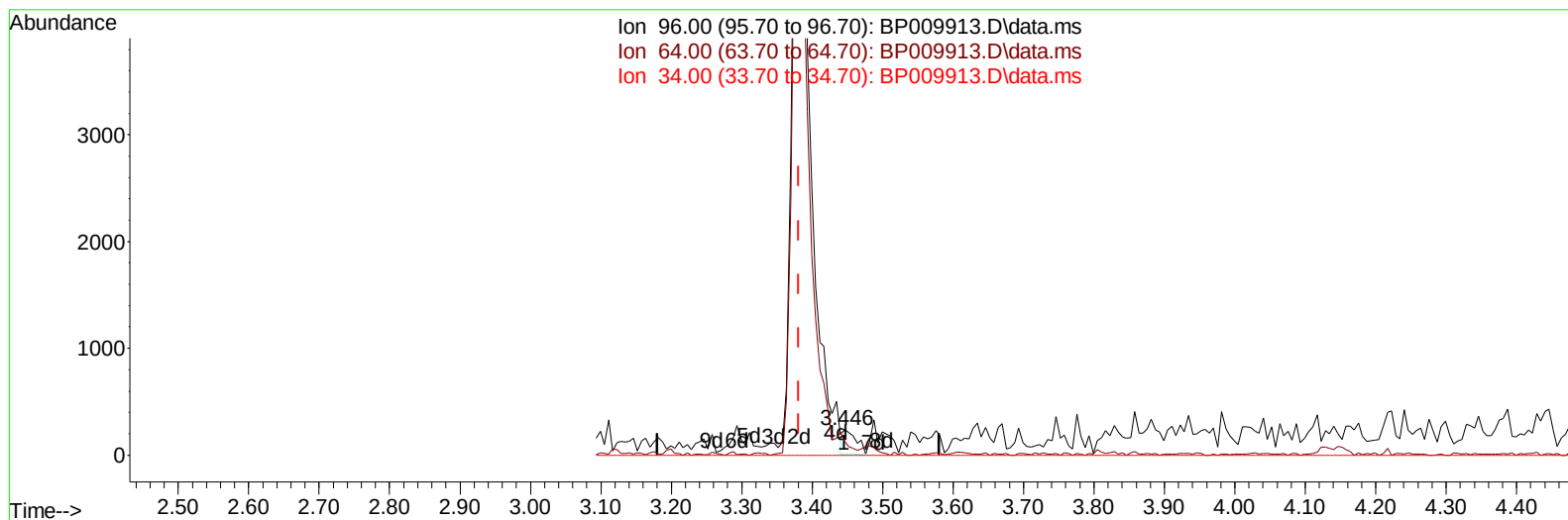
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009913.D
 Acq On : 18 Apr 2022 13:53
 Operator : CG/JU
 Sample : N2422-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J03

Manual IntegrationsAPPROVED

Quant Time: Apr 19 00:55:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009913.D\data.ms

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

3.446min (+ 0.065) 0.12 ng/uL

response 295

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	54.20
34.00	0.00	0.00
0.00	0.00	0.00

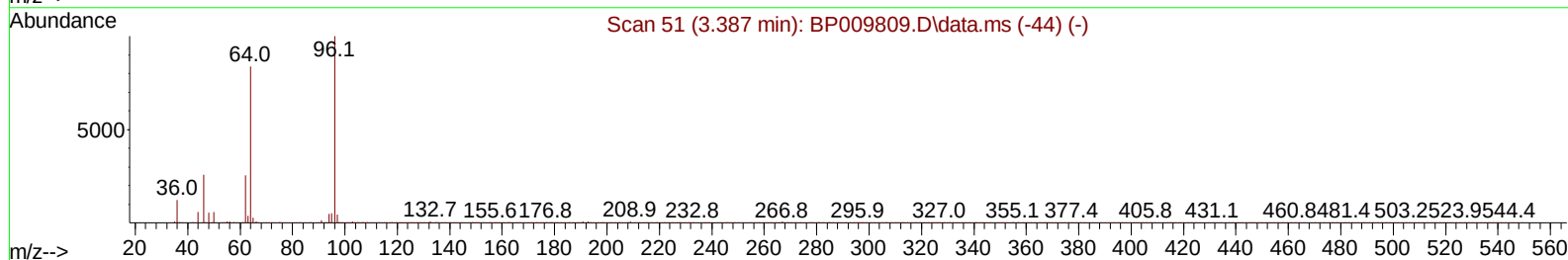
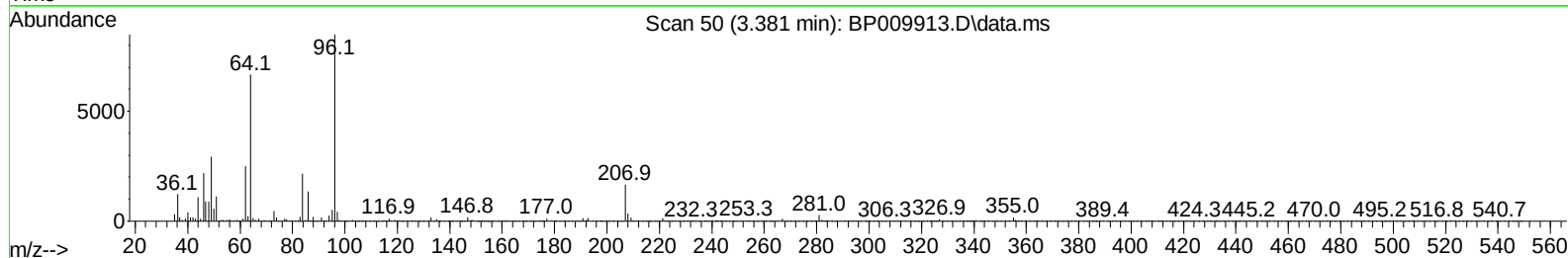
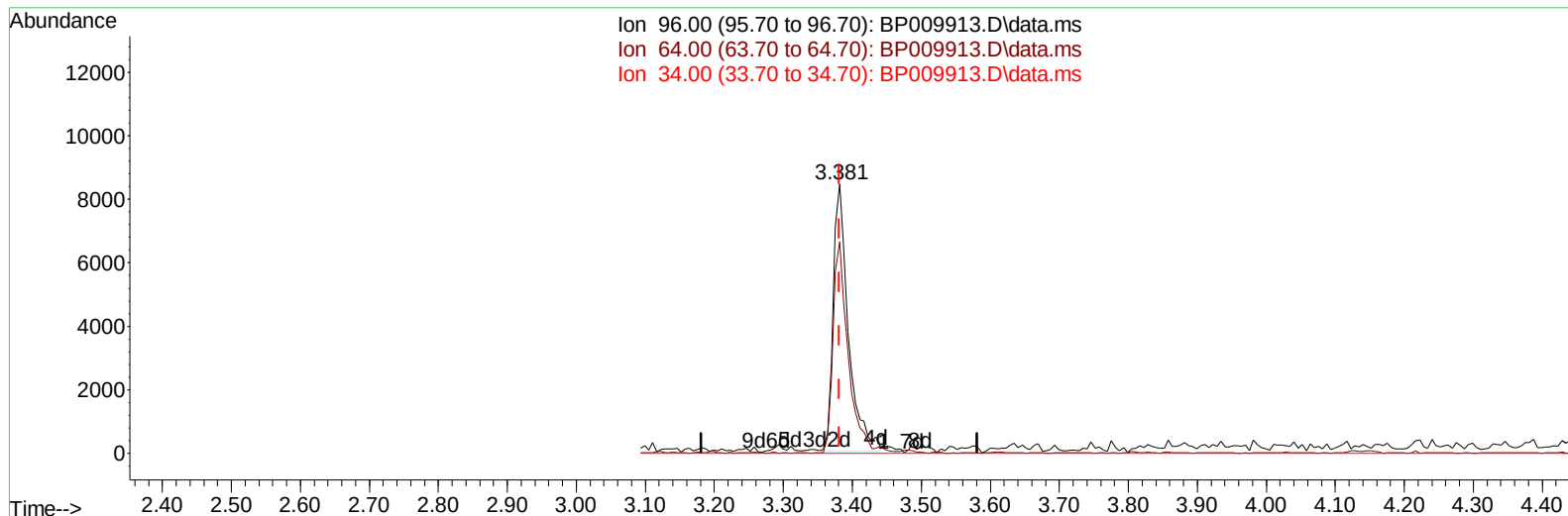
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(3) 1,4-Dioxane-d8 (S)

3.381min (+ 0.000) 5.03 ng/uL m

response 12861

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.58#
34.00	0.00	0.00
0.00	0.00	0.00

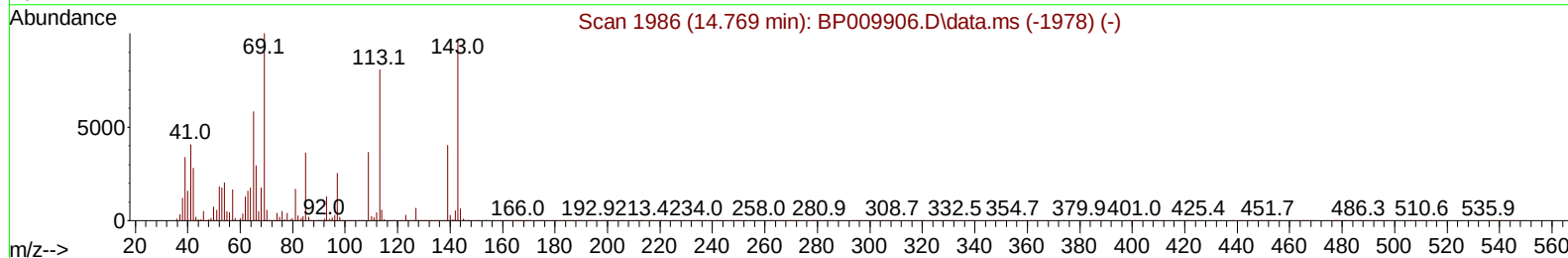
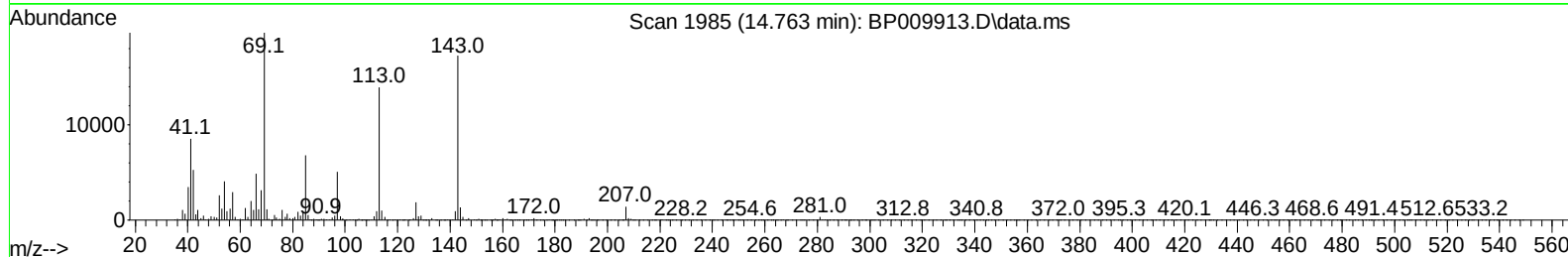
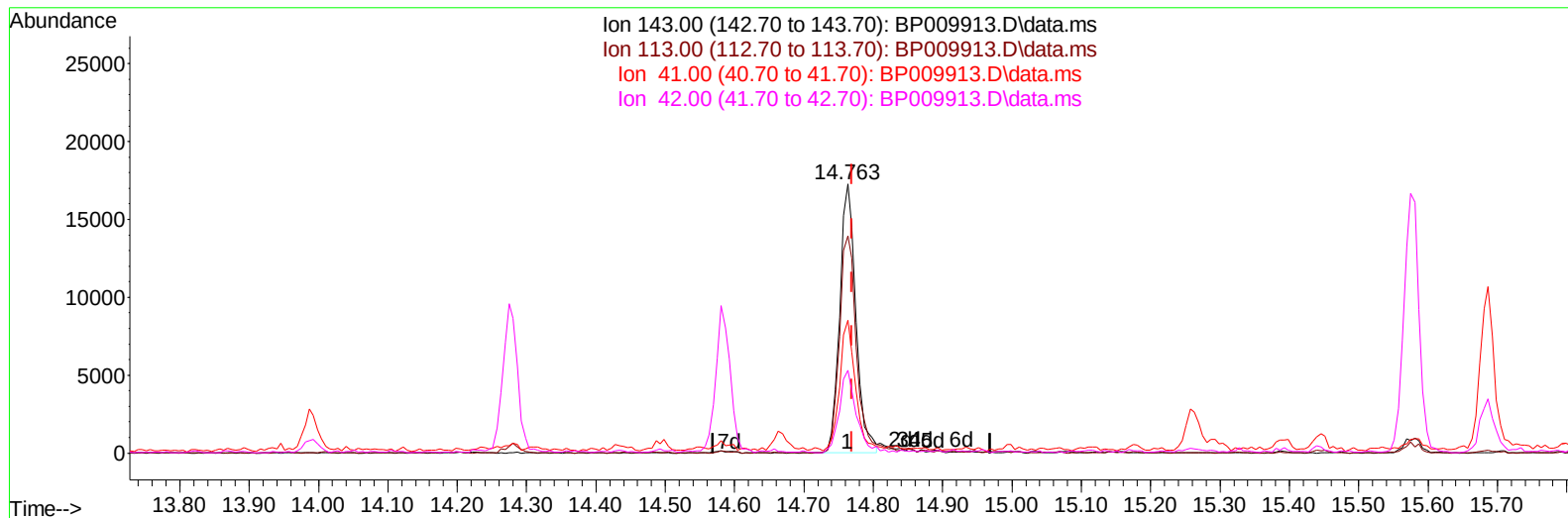
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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TIC: BP009913.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.03 ng/ul

response 27685

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.72#
41.00	23.20	49.47#
42.00	18.00	30.67#

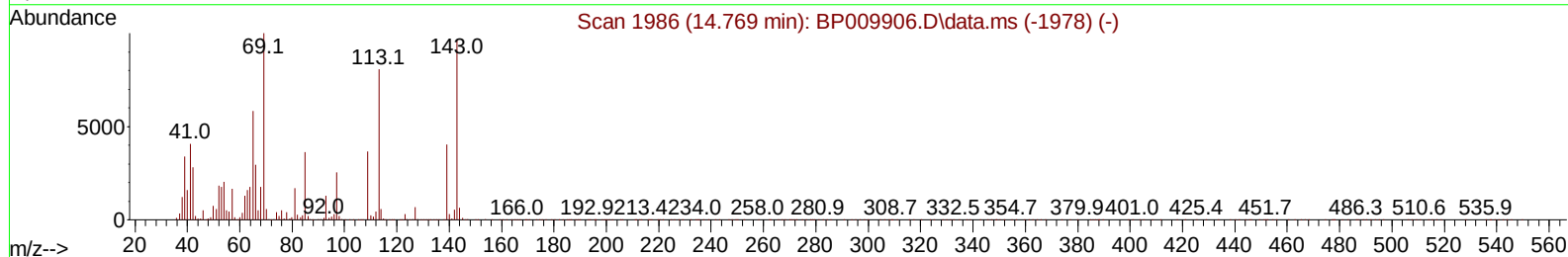
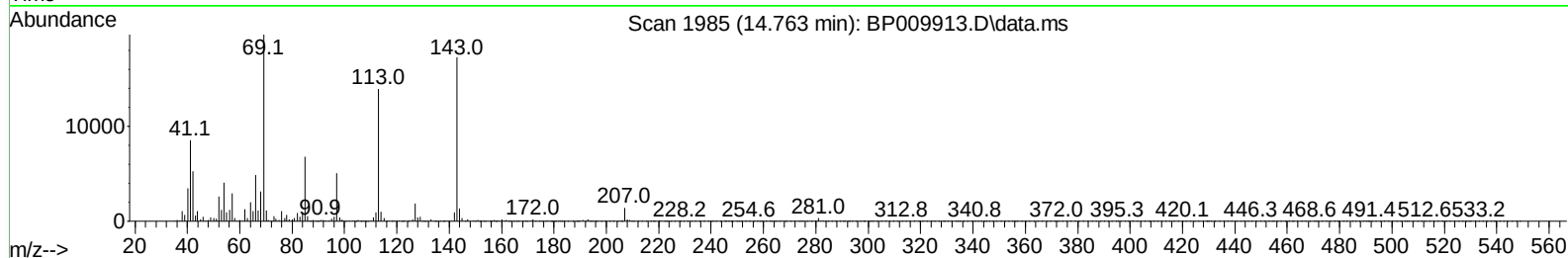
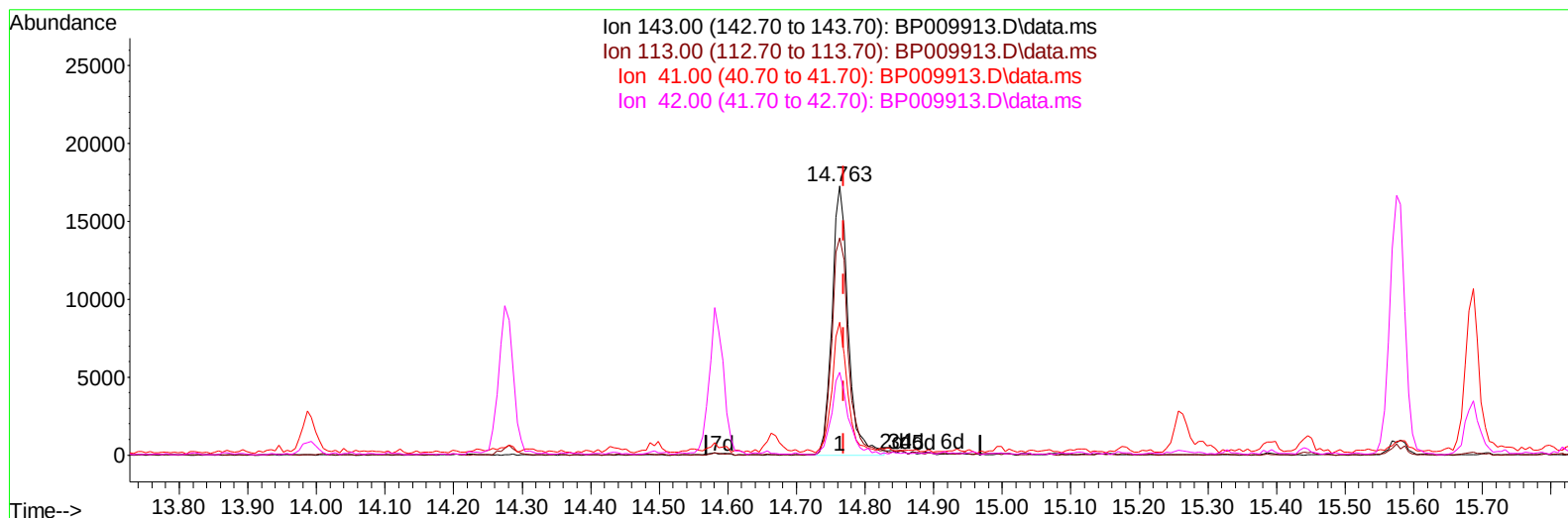
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Operator : CG/JU
Sample : N2422-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP009913.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 5.16 ng/ul m

response 28372

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.72#
41.00	23.20	49.47#
42.00	18.00	30.67#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	114054	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.757	136	513484	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.587	164	386320	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.334	188	873226	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	925970	20.000	ng/ul	# 0.00
88) Perylene-d12	23.874	264	851080	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	12861m	5.033	ng/uL	0.00
4) Pyridine-d5	3.805	84	41669	5.914	ng/ul	0.00
7) Phenol-d5	7.111	99	56143	5.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	186463	31.279	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	169054	22.382	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	119835	14.336	ng/ul	-0.01
21) Nitrobenzene-d5	9.111	128	119582	32.294	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	124772	30.735	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	220753	25.986	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	302839	25.029	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1001976	33.018	ng/ul	0.00
49) Acenaphthylene-d8	14.275	160	1069987	29.762	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	28372m	5.155	ng/ul	0.00
60) Fluorene-d10	15.575	176	876068	34.318	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	155273	30.016	ng/ul	0.00
73) Anthracene-d10	17.434	188	1557424	37.251	ng/ul	0.00
81) Pyrene-d10	19.663	212	1936121	38.284	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	1709472	38.570	ng/ul	-0.01
Target Compounds						
35) 4-Chloro-3-methylphenol	12.034	107	20546	2.059	ng/ul#	Qvalue 71

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

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