

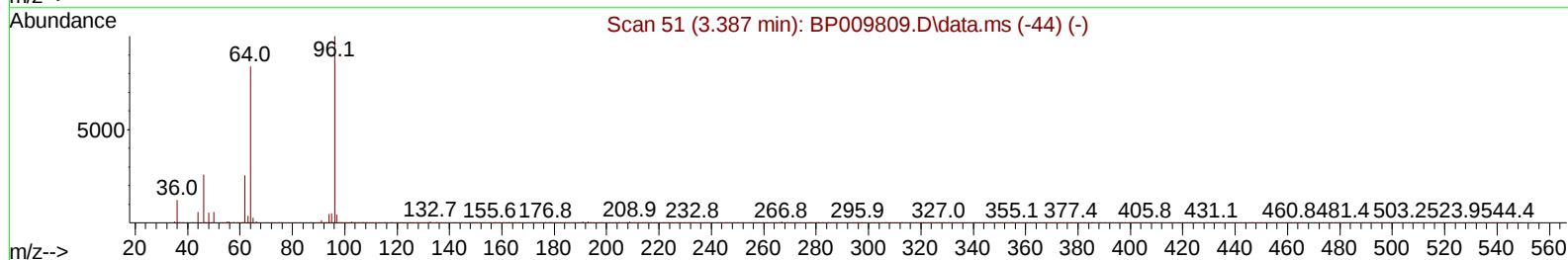
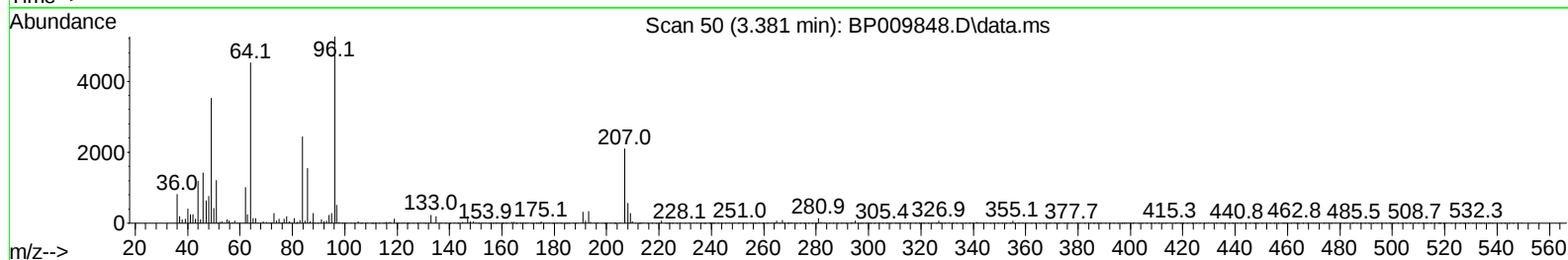
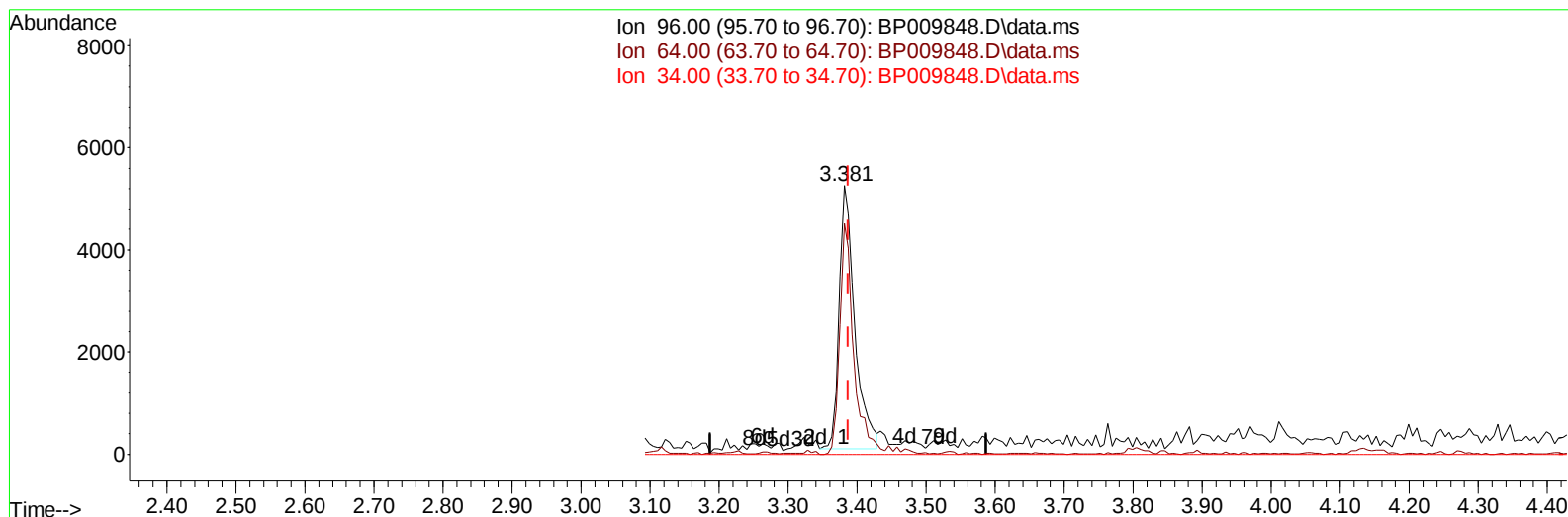
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009848.D
 Acq On : 15 Apr 2022 03:06
 Operator : CG/JU
 Sample : N2293-19
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNP7

Manual IntegrationsAPPROVED

Quant Time: Apr 15 04:22:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 13 02:15:22 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/22/2022



TIC: BP009848.D\data.ms

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

3.381min (-0.006) 2.63 ng/uL

response 8139

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.04#
34.00	0.00	0.00
0.00	0.00	0.00

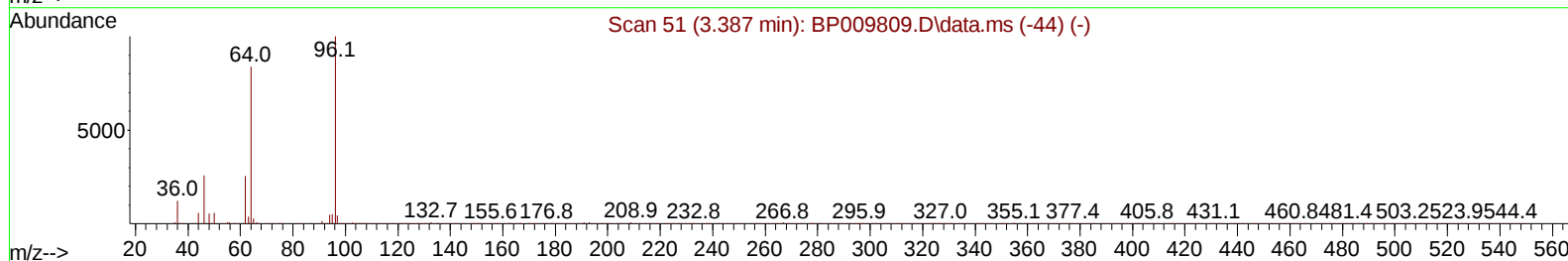
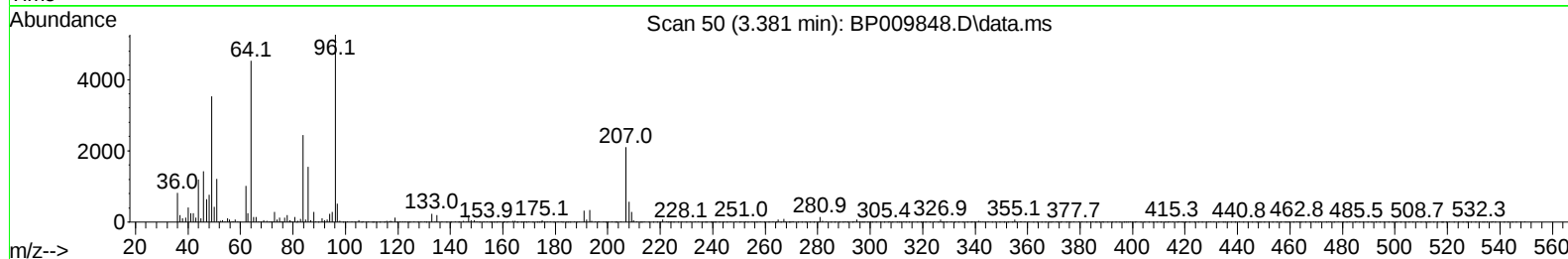
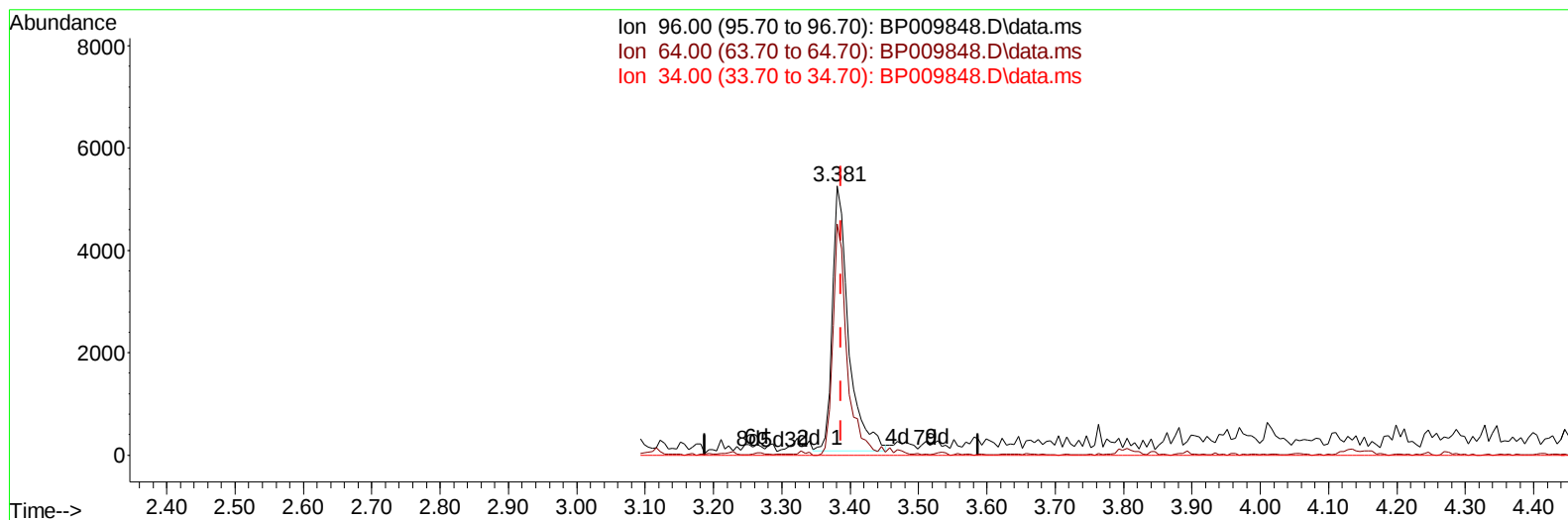
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TIC: BP009848.D\data.ms

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

3.381min (-0.006) 2.76 ng/uL m

response 8532

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.04#
34.00	0.00	0.00
0.00	0.00	0.00

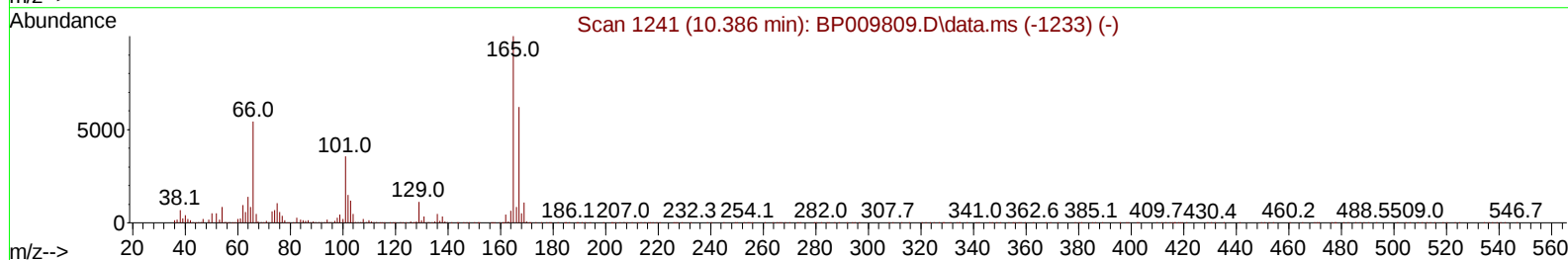
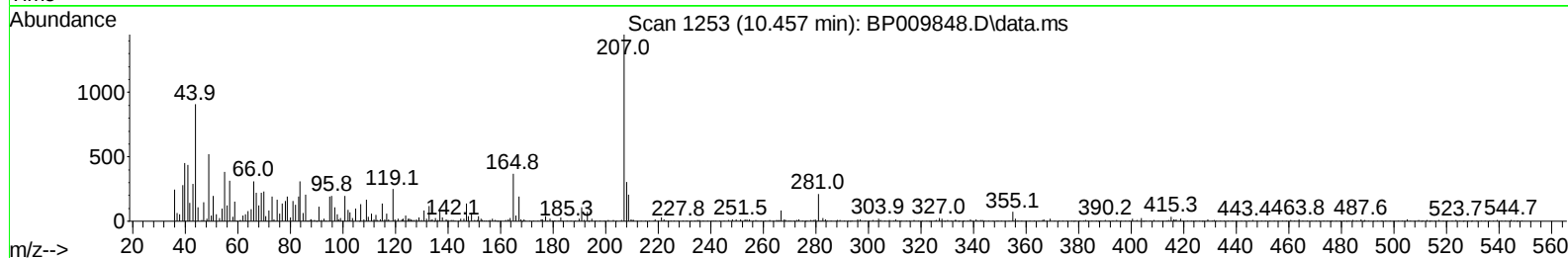
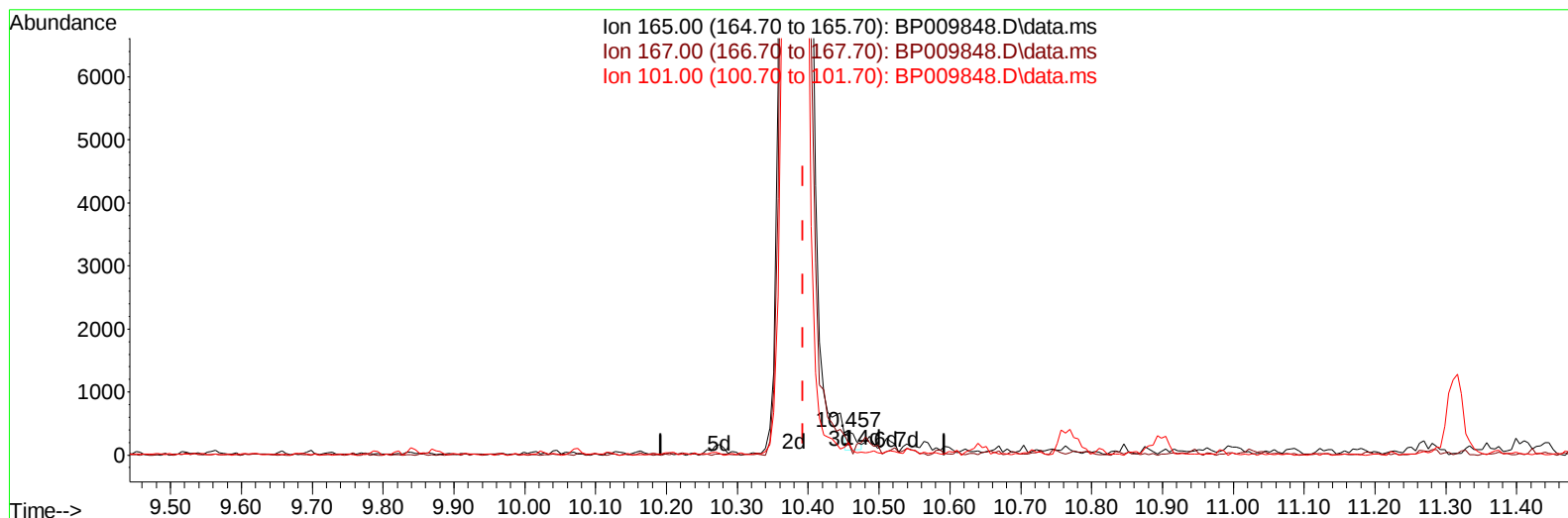
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TIC: BP009848.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.457min (+ 0.065) 0.03 ng/u1

response 272

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	51.76
101.00	49.80	53.93
0.00	0.00	0.00

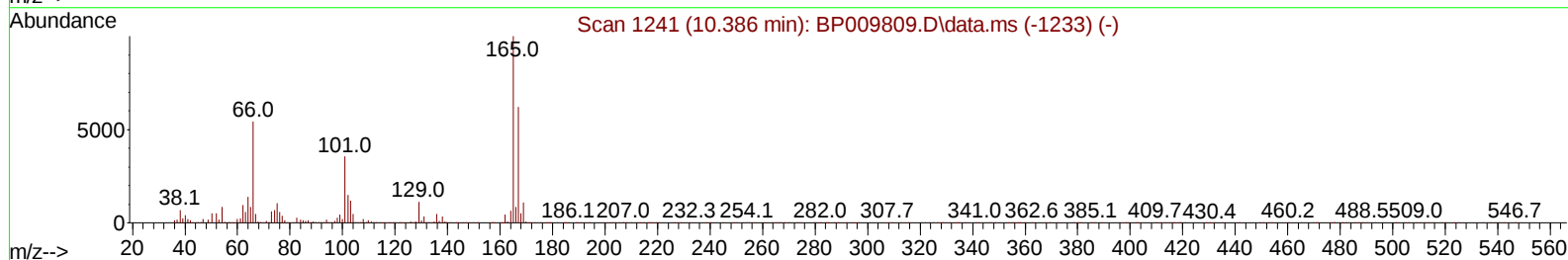
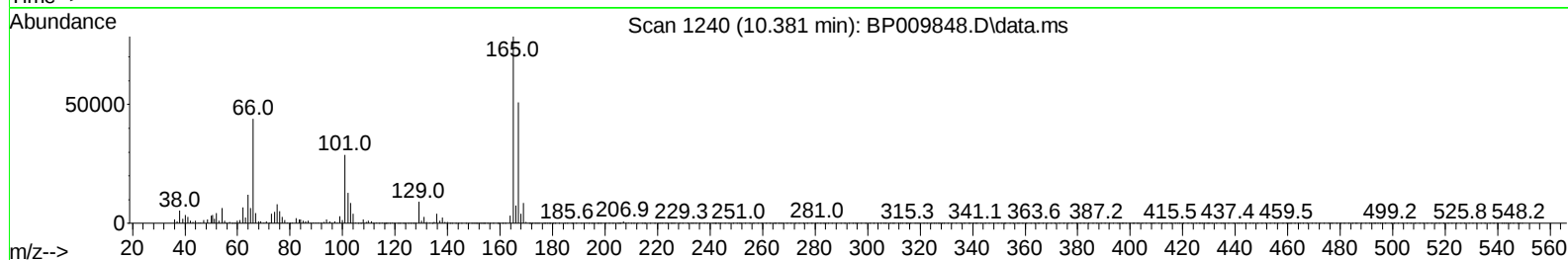
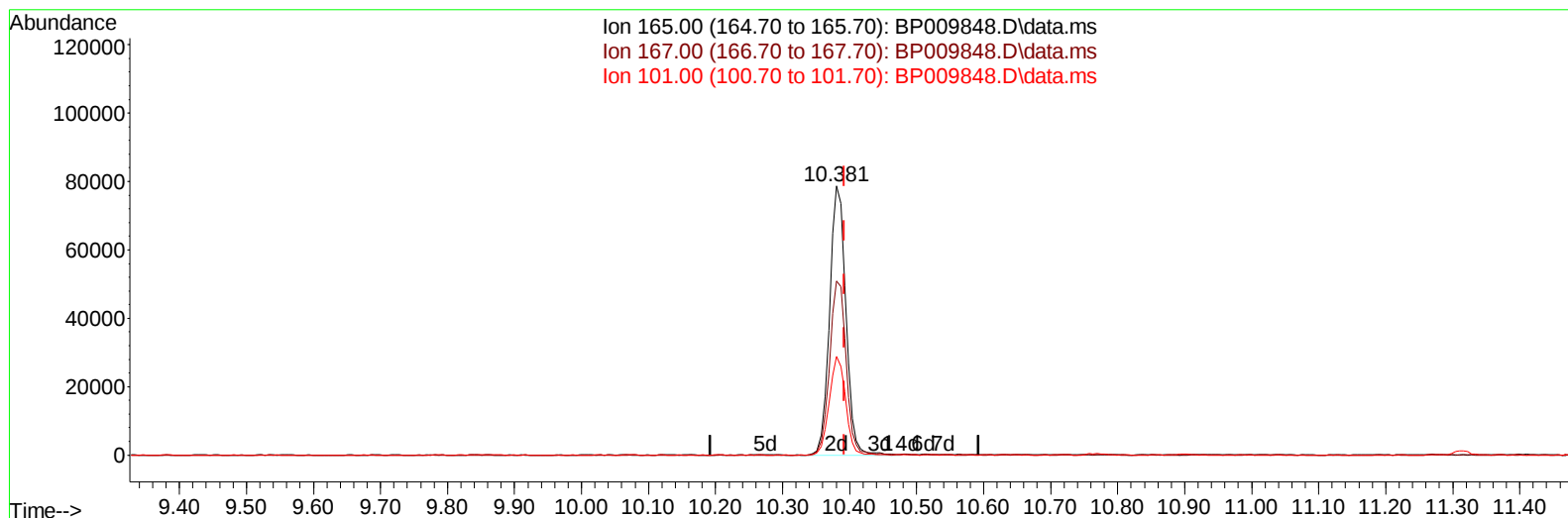
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Instrument :
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TIC: BP009848.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.381min (-0.012) 13.03 ng/ul m

response 132757

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	64.70
101.00	49.80	36.78#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.963	152	138051	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	616070	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	432863	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	969192	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.421	240	1035604	20.000	ng/ul	#-0.01
88) Perylene-d12	23.880	264	1021453	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	8532m	2.758	ng/uL	0.00
4) Pyridine-d5	3.805	84	71443	8.377	ng/ul	0.00
7) Phenol-d5	7.110	99	159149	13.131	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.281	67	110748	15.349	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.487	132	130371	14.260	ng/ul	-0.01
15) 4-Methylphenol-d8	8.657	113	122796	12.137	ng/ul	-0.01
21) Nitrobenzene-d5	9.116	128	70755	15.926	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	73331	15.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	132757m	13.025	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.898	131	169190	11.655	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	560085	16.472	ng/ul	-0.01
49) Acenaphthylene-d8	14.281	160	569597	14.140	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	78955	12.804	ng/ul	-0.01
60) Fluorene-d10	15.581	176	490733	17.156	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	54179	9.436	ng/ul	0.00
73) Anthracene-d10	17.439	188	836452	18.025	ng/ul	-0.01
81) Pyrene-d10	19.669	212	1145863	20.259	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1024884	19.267	ng/ul	-0.02

Target Compounds Qvalue

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

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