

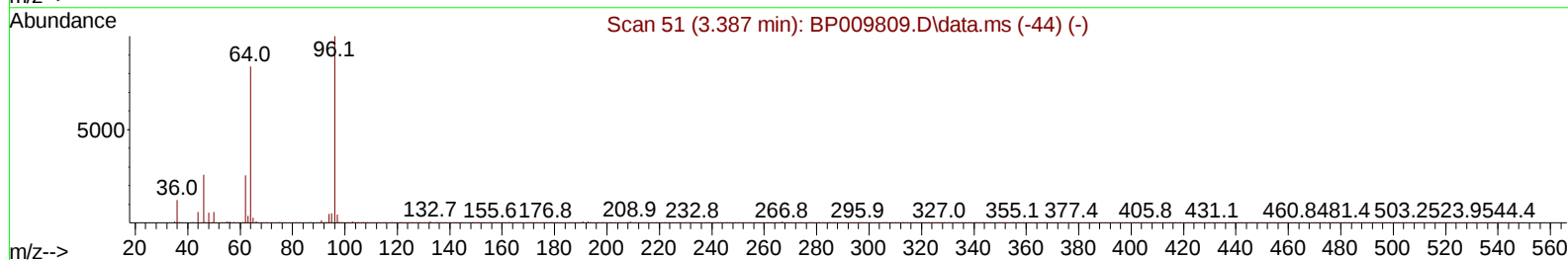
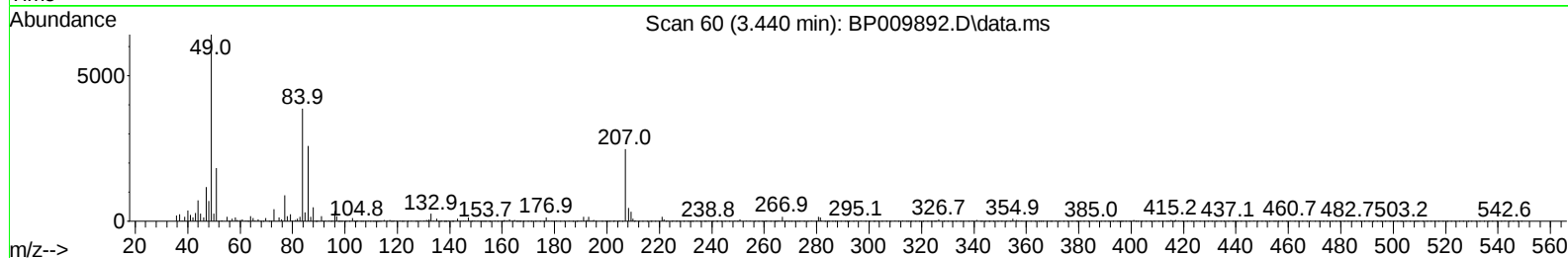
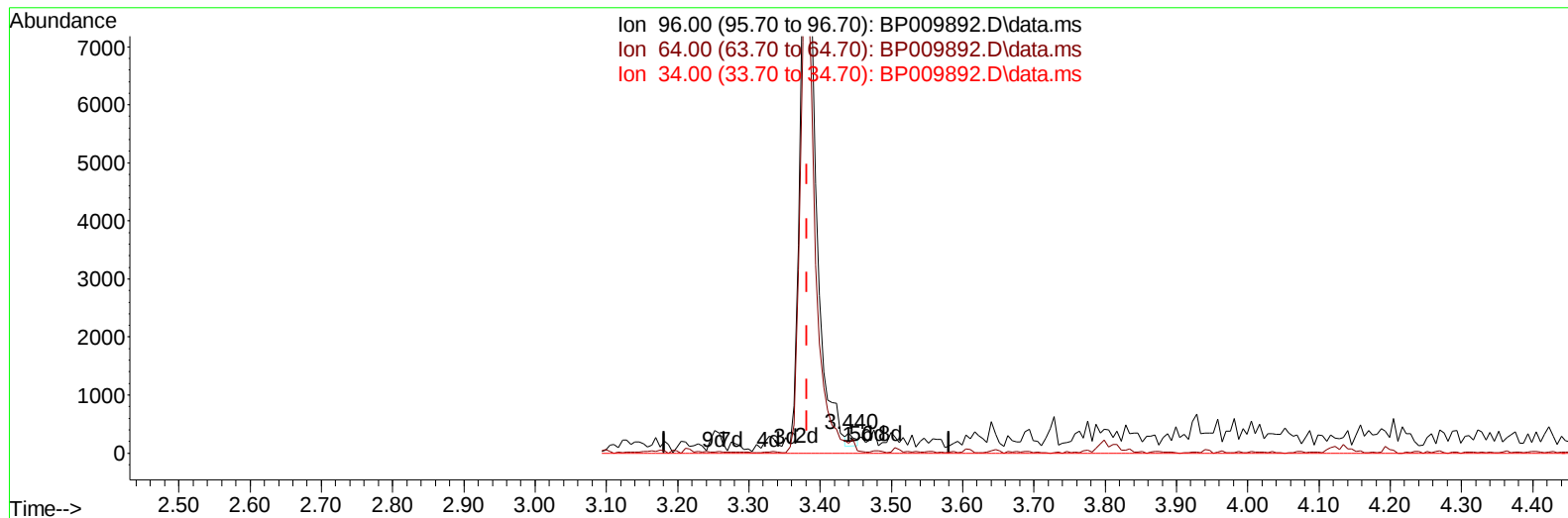
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
Data File : BP009892.D
Acq On : 16 Apr 2022 09:34
Operator : CG/JU
Sample : N2421-10
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HY7

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:10:55 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 04/18/2022
Supervised By :Yogesh Patel 04/18/2022



TIC: BP009892.D\data.ms

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

3.440min (+ 0.059) 0.05 ng/uL

response 164

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	51.61
34.00	0.00	0.00
0.00	0.00	0.00

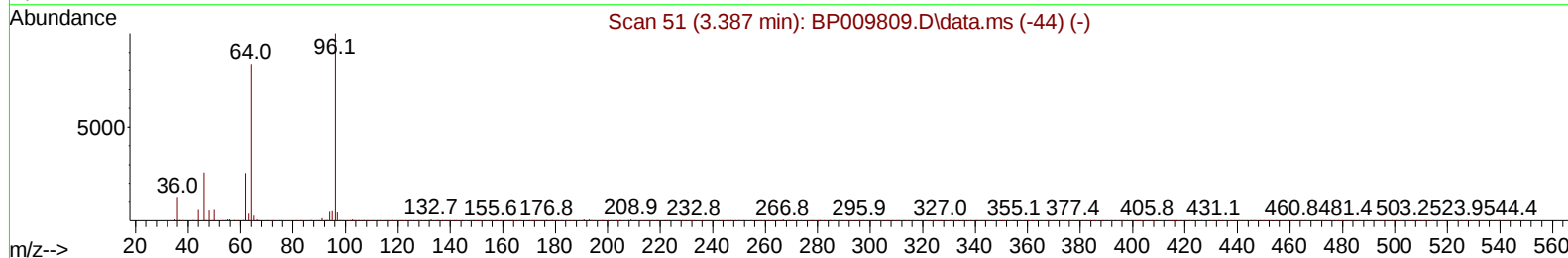
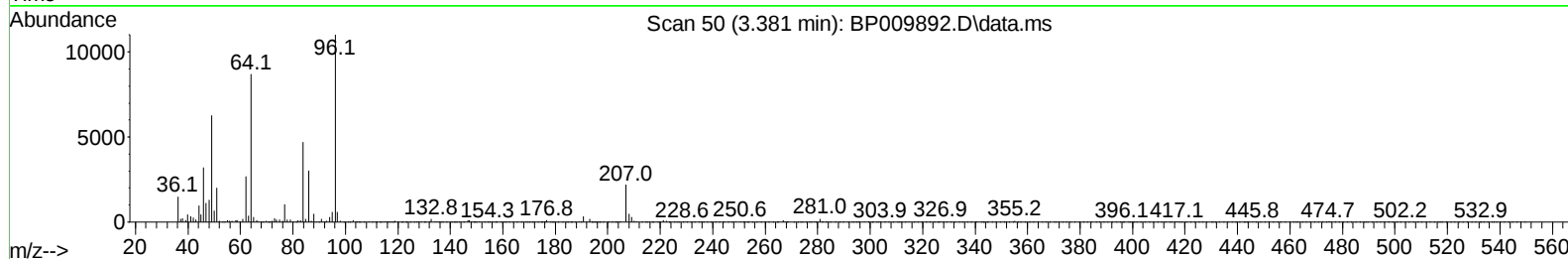
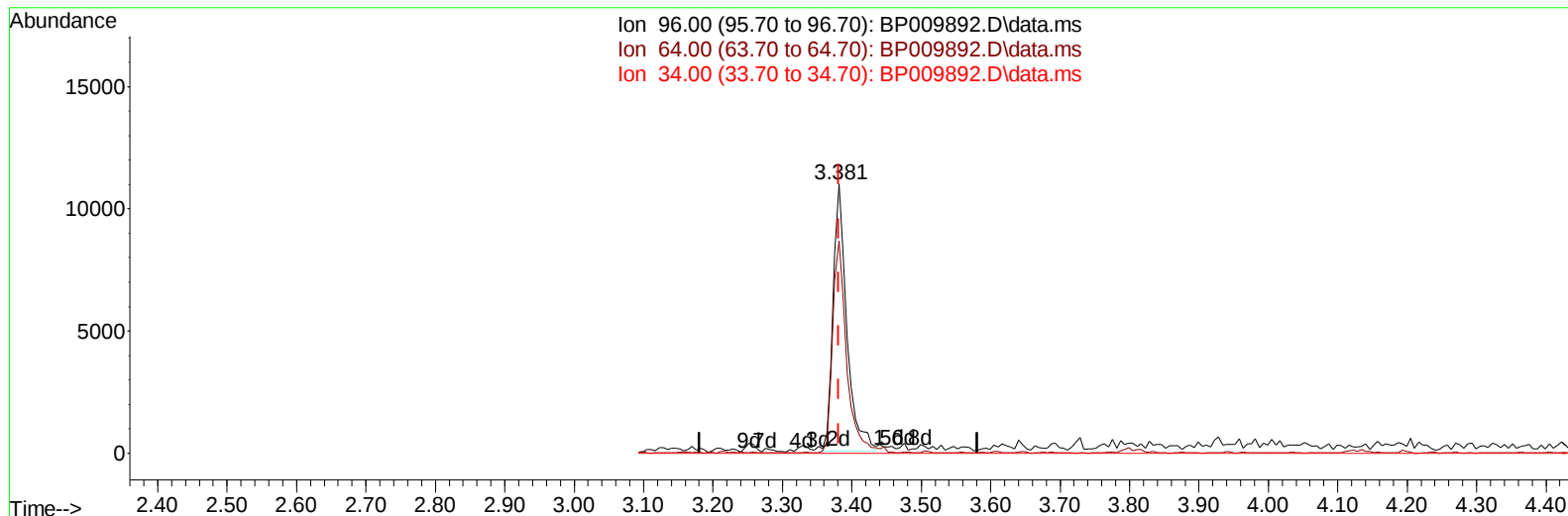
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TIC: BP009892.D\data.ms

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

3.381min (+ 0.000) 4.86 ng/uL m

response 15126

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.87#
34.00	0.00	0.00
0.00	0.00	0.00

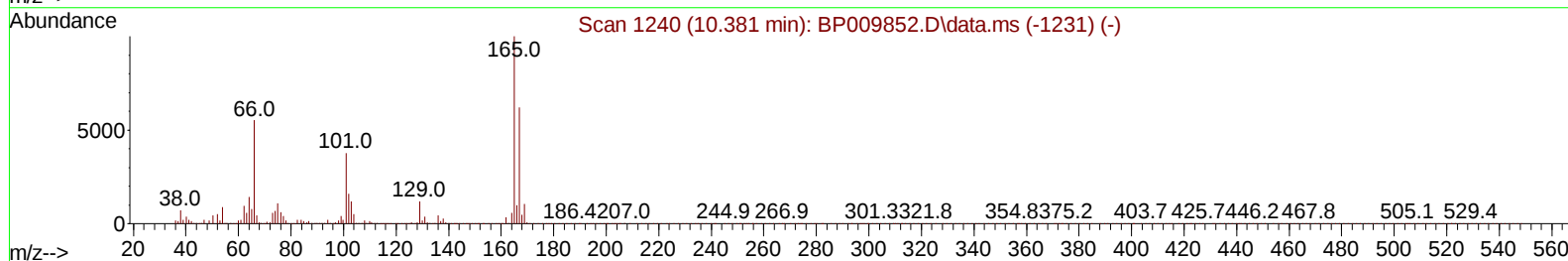
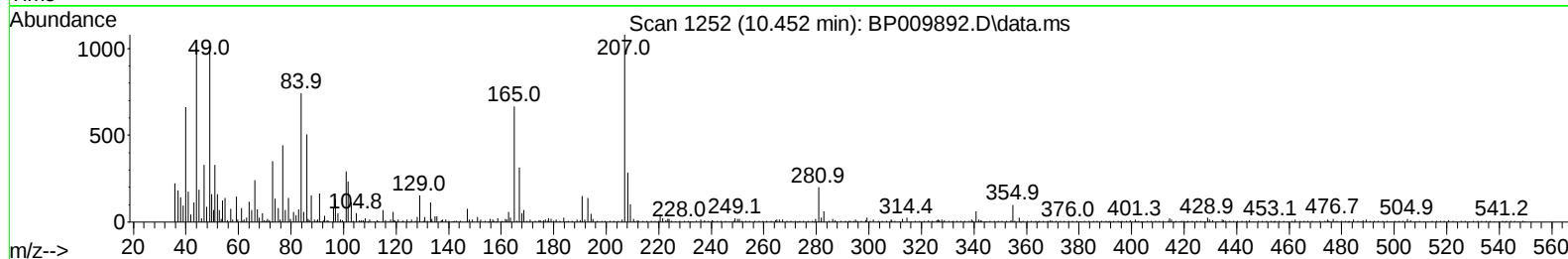
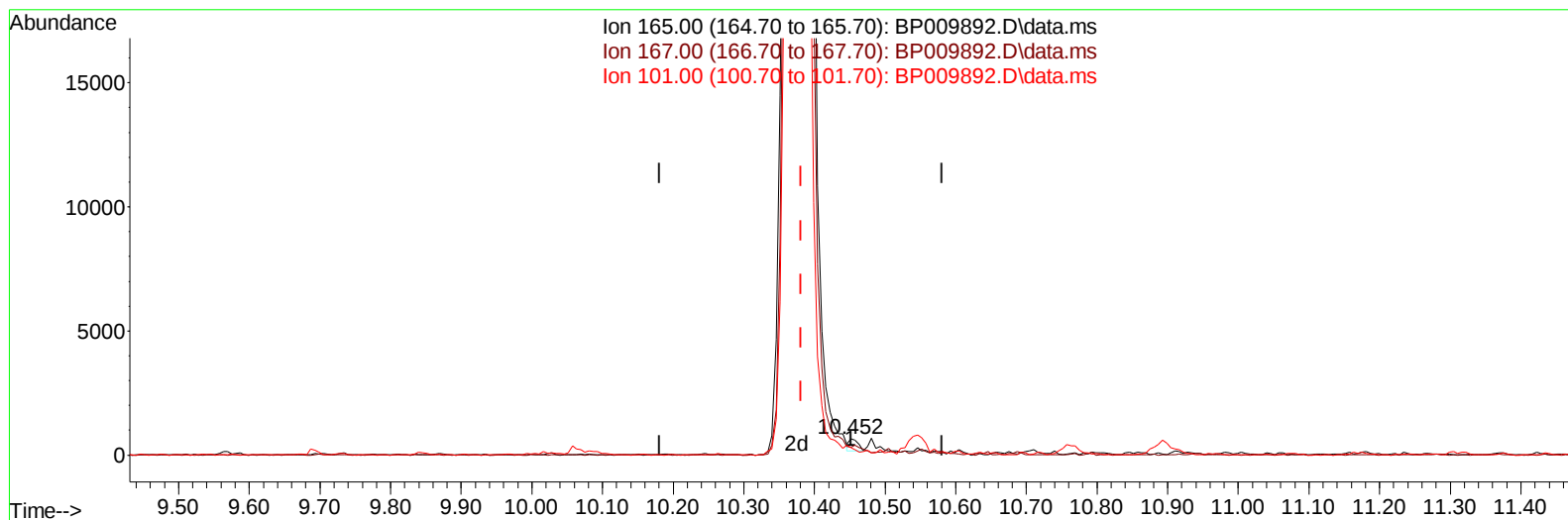
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TIC: BP009892.D\data.ms

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

10.452min (+ 0.071) 0.04 ng/ul

response 408

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	47.45
101.00	49.80	43.69
0.00	0.00	0.00

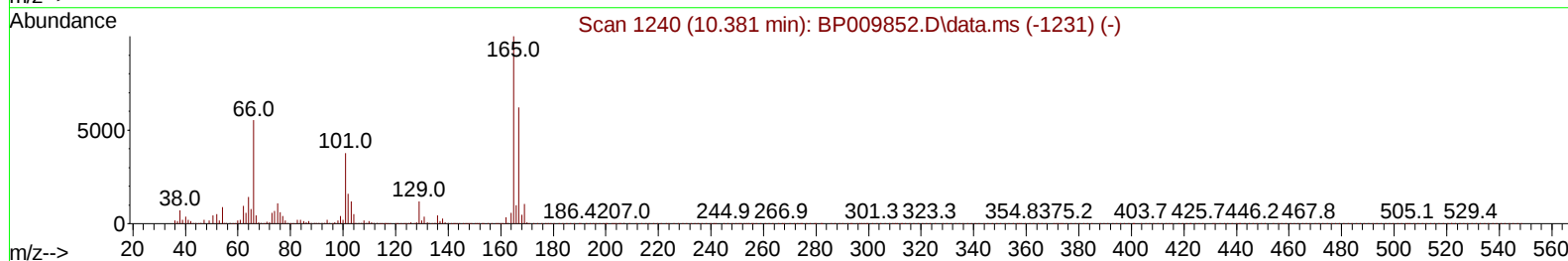
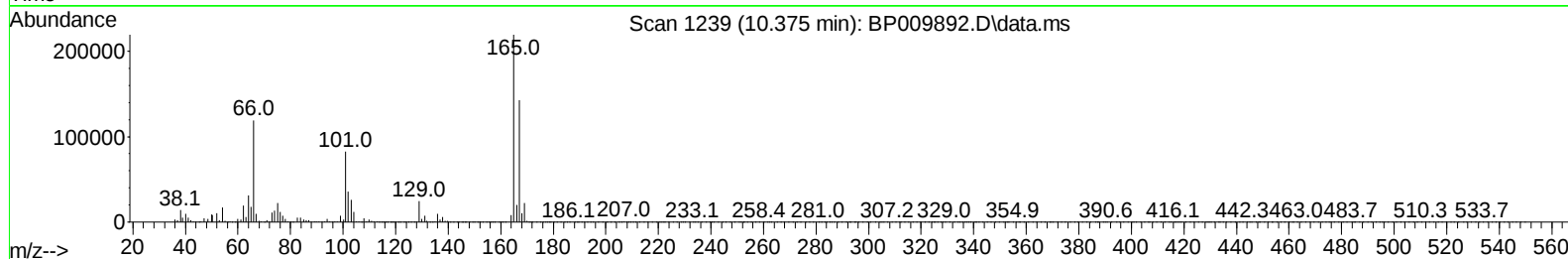
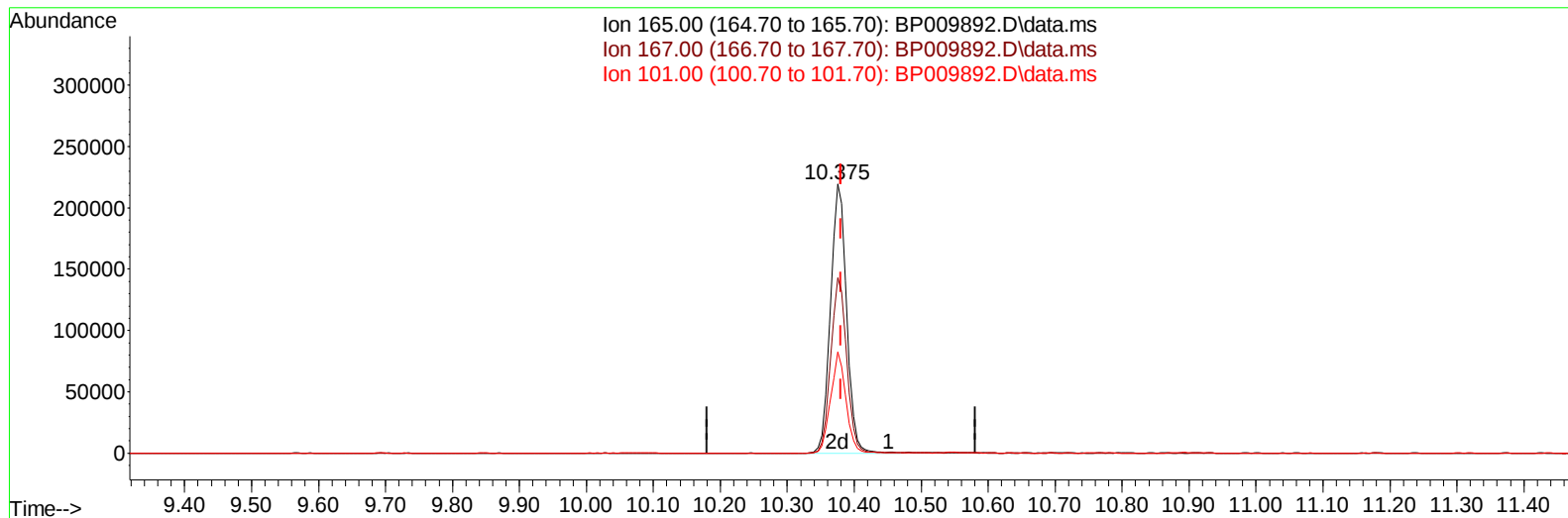
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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP009892.D\data.ms

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

10.375min (-0.006) 35.57 ng/ul m

response 366461

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	65.22
101.00	49.80	37.72#
0.00	0.00	0.00

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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	7.958	152	138937	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	622670	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	442694	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	960951	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	982665	20.000	ng/ul	# 0.00
88) Perylene-d12	23.880	264	755973	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	15126m	4.859	ng/uL	0.00
4) Pyridine-d5	3.799	84	93749	10.922	ng/ul	0.00
7) Phenol-d5	7.111	99	108407	8.887	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	268738	37.007	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	272545	29.621	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	198963	19.540	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	178670	39.791	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	194019	39.412	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	366461m	35.574	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	294595	20.078	ng/ul	0.00
46) Dimethylphthalate-d6	13.993	166	1409308	40.527	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1533515	37.223	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	55004	8.722	ng/ul	0.00
60) Fluorene-d10	15.581	176	1212180	41.437	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	216815	38.086	ng/ul	0.00
73) Anthracene-d10	17.439	188	1980669	43.049	ng/ul	0.00
81) Pyrene-d10	19.669	212	2436155	45.393	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1857840	47.191	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.134	94	19265	1.578	ng/ul#	85
86) Bis(2-ethylhexyl)phtha...	21.327	149	49898	1.291	ng/ul#	83

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

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