

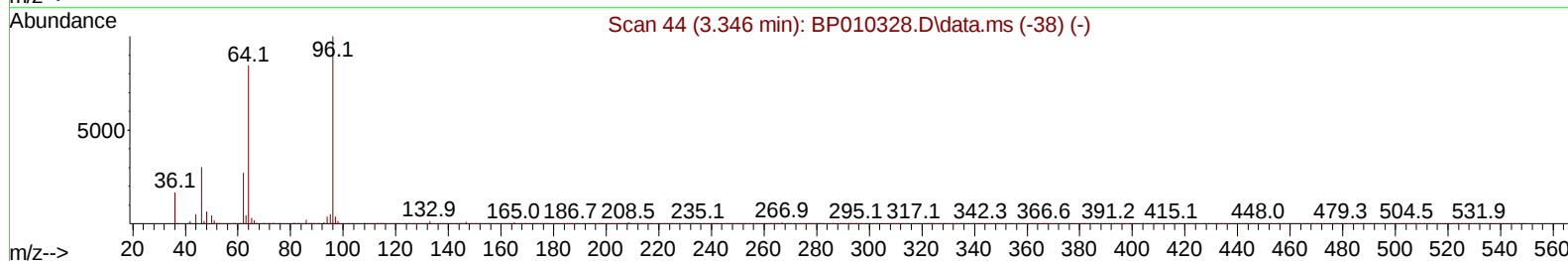
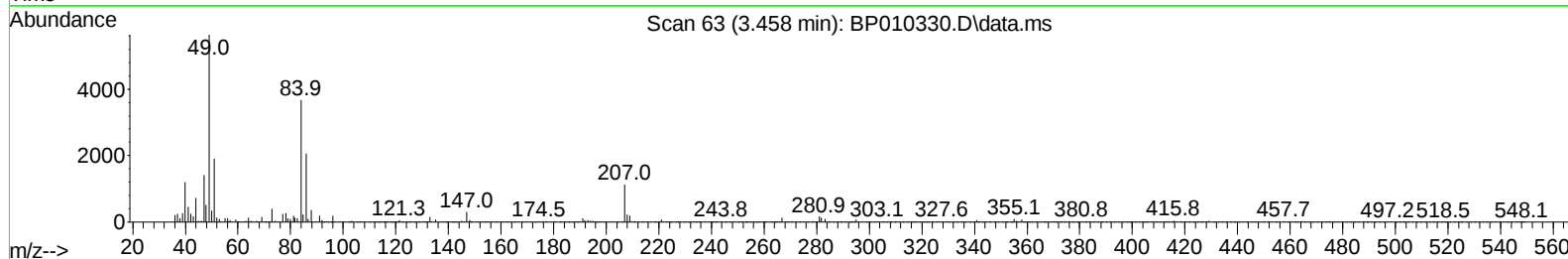
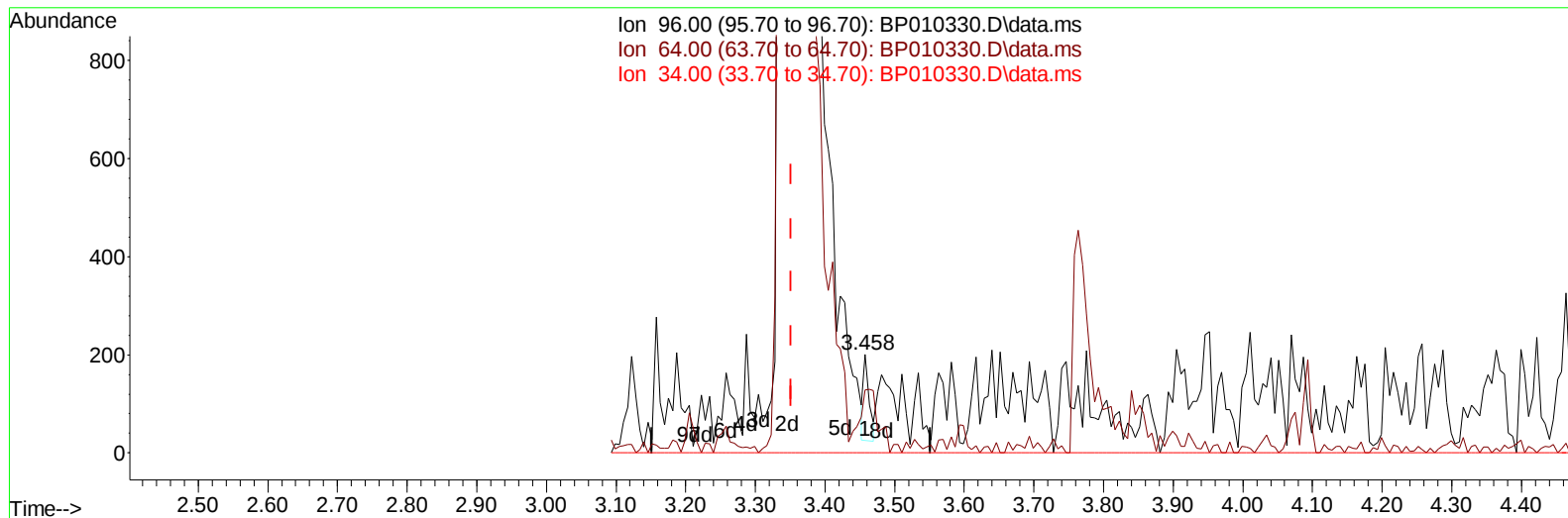
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010330.D
 Acq On : 12 May 2022 19:08
 Operator : CG/JU
 Sample : PB144782BL
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK782

Manual IntegrationsAPPROVED

Quant Time: May 13 01:15:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/16/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010330.D\data.ms

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

3.458min (+ 0.106) 0.03 ng/uL

response 101

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	64.18
34.00	0.00	0.00
0.00	0.00	0.00

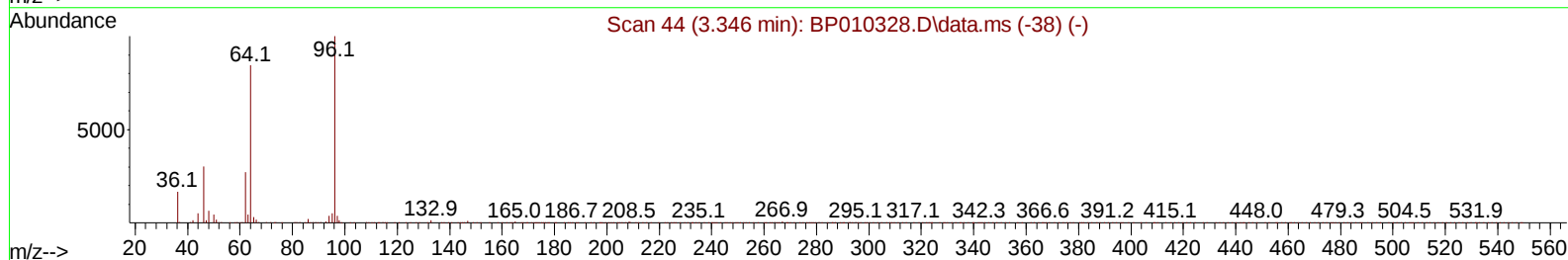
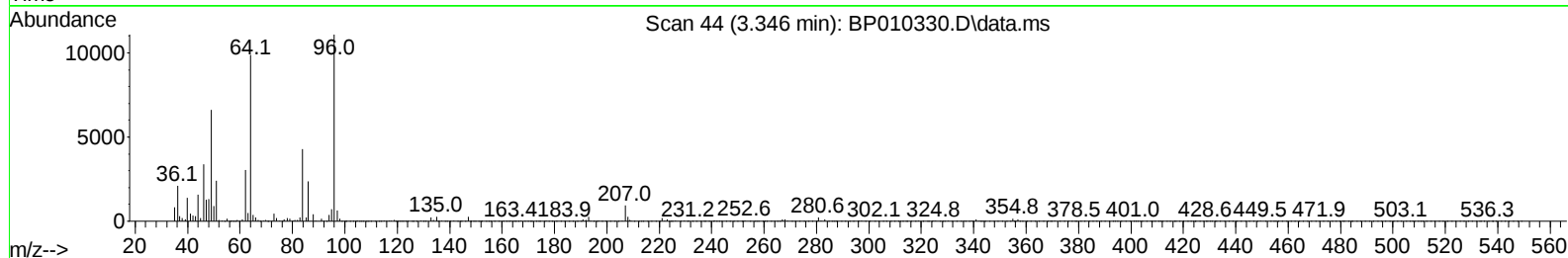
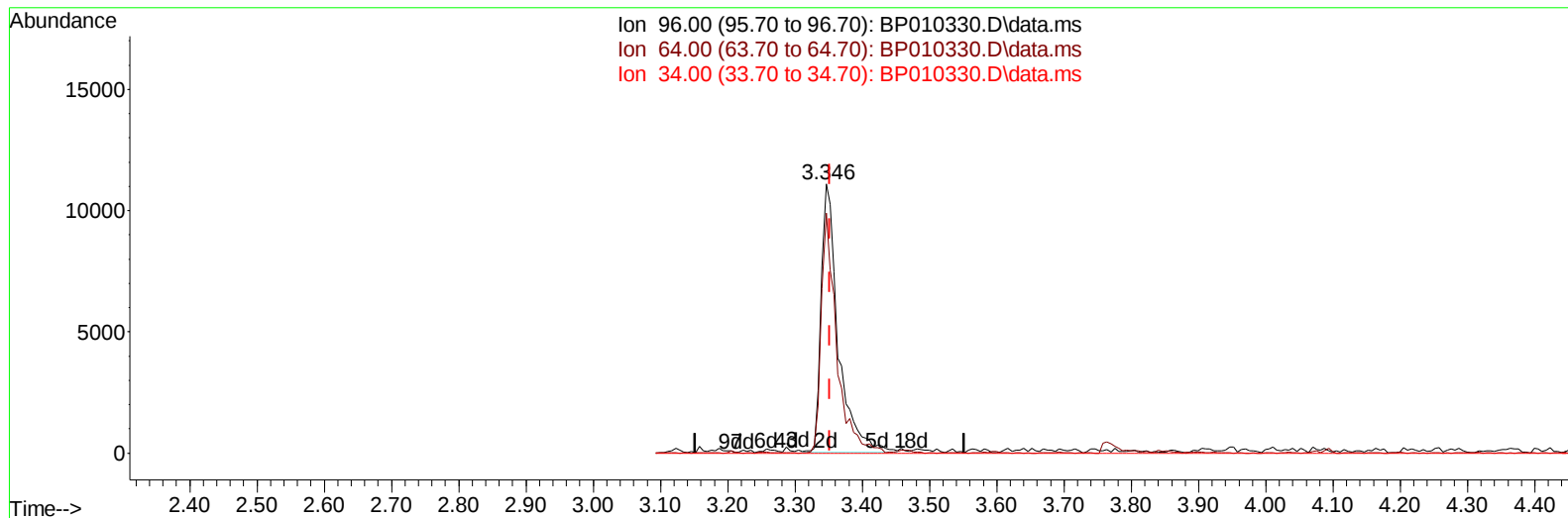
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TIC: BP010330.D\data.ms

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

3.346min (-0.006) 6.41 ng/uL m

response 19580

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.23#
34.00	0.00	0.00
0.00	0.00	0.00

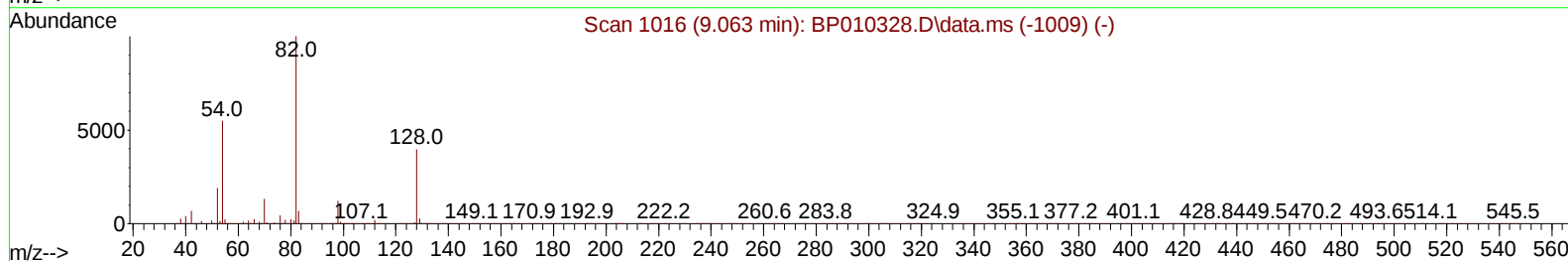
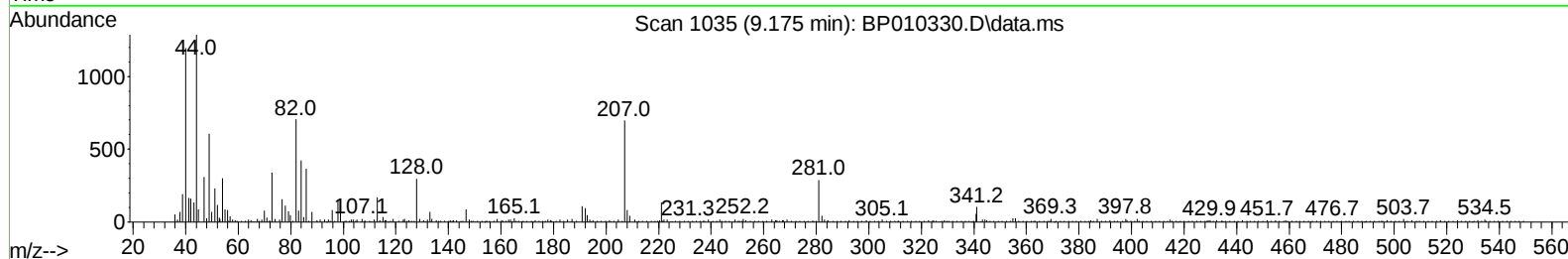
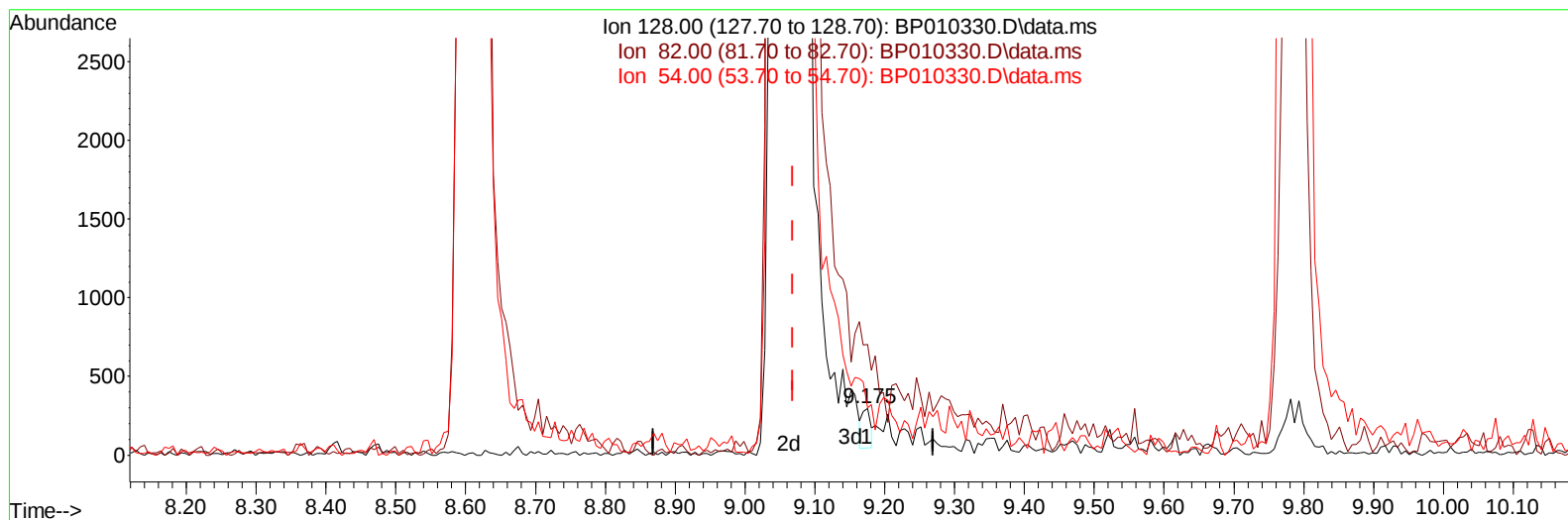
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Misc :
ALS Vial : 54 Sample Multiplier: 1

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TIC: BP010330.D\data.ms

(21) Nitrobenzene-d5 (S)

9.175min (+ 0.106) 0.05 ng/u1

response

215

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	236.58
54.00	120.20	100.67
0.00	0.00	0.00

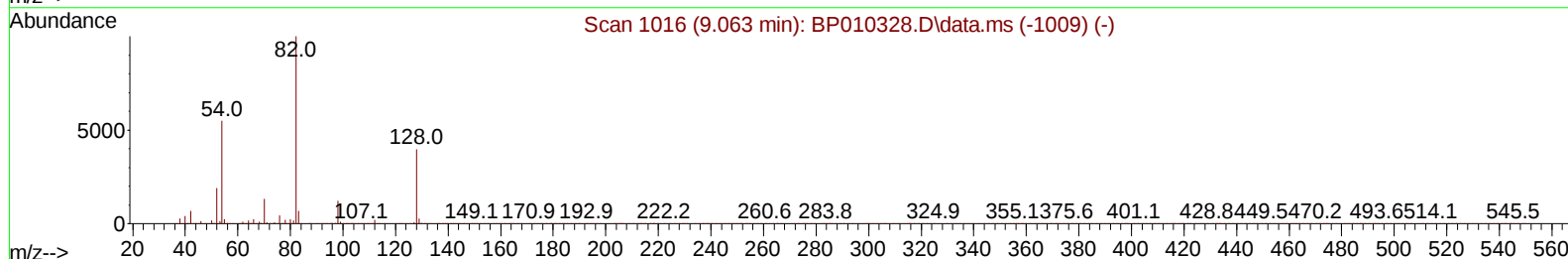
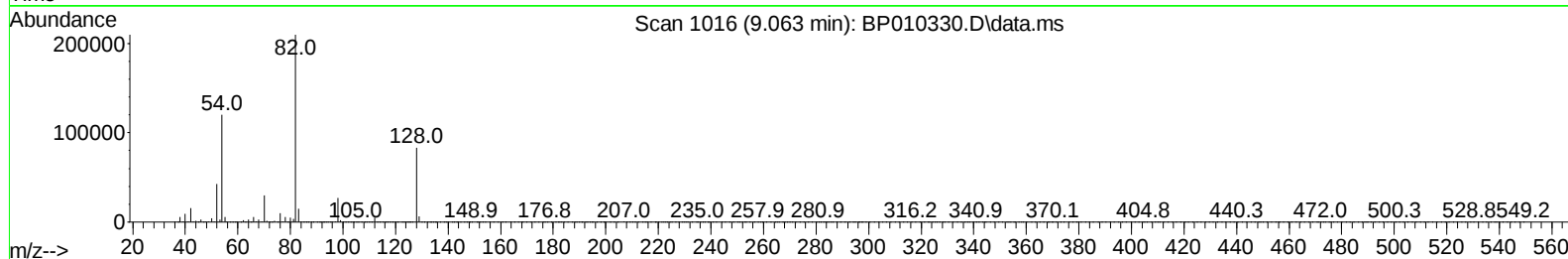
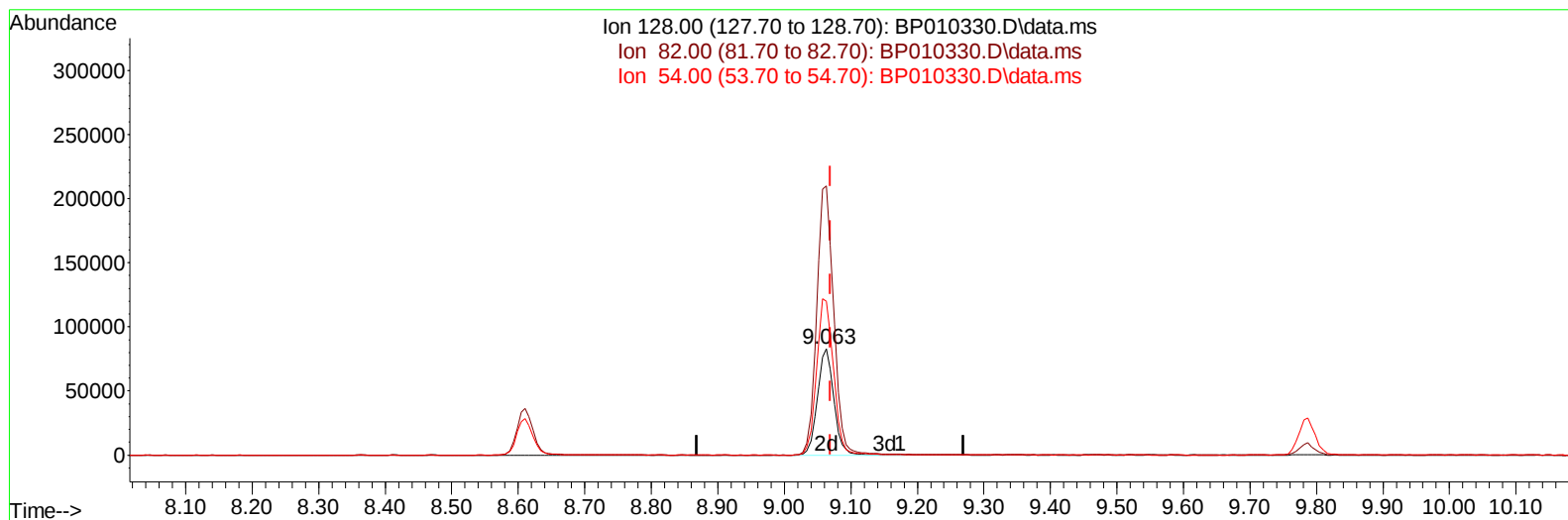
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Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP010330.D\data.ms

(21) Nitrobenzene-d5 (S)

9.063min (-0.006) 31.44 ng/ul m

response 142586

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	252.75
54.00	120.20	144.82#
0.00	0.00	0.00

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 Misc :
 ALS Vial : 54 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	129698	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	565601	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	403958	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	892360	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	905604	20.000	ng/ul	-0.01
88) Perylene-d12	23.792	264	847601	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	19580m	6.413	ng/uL	0.00
4) Pyridine-d5	3.764	84	262617	30.911	ng/ul	0.00
7) Phenol-d5	7.069	99	325564	27.026	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	230964	30.993	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	262576	29.597	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	283874	28.664	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	142586m	31.441	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	155069	29.990	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	264721	28.128	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	405463	29.183	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1014964	31.538	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1133831	29.551	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	145202	23.522	ng/ul	-0.01
60) Fluorene-d10	15.528	176	861672	31.719	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	138452	22.657	ng/ul	-0.01
73) Anthracene-d10	17.386	188	1390016	31.809	ng/ul	0.00
81) Pyrene-d10	19.616	212	1685315	33.676	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1480480	32.752	ng/ul	-0.02

Target Compounds Qvalue

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

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