

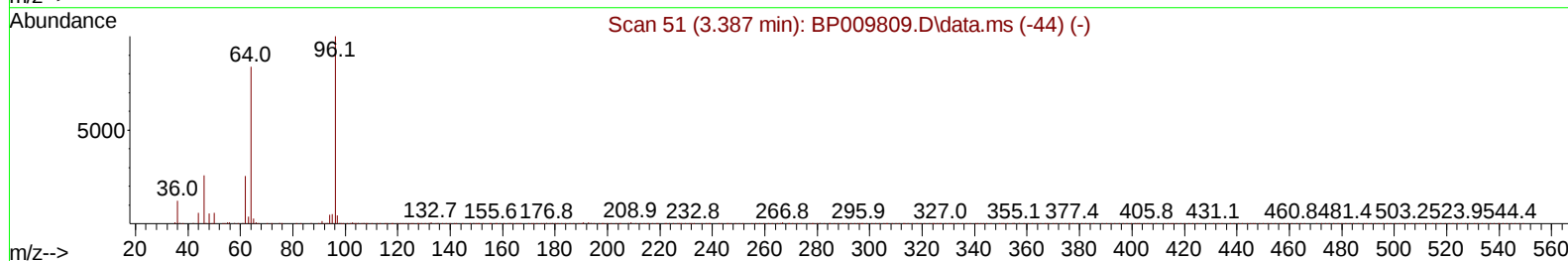
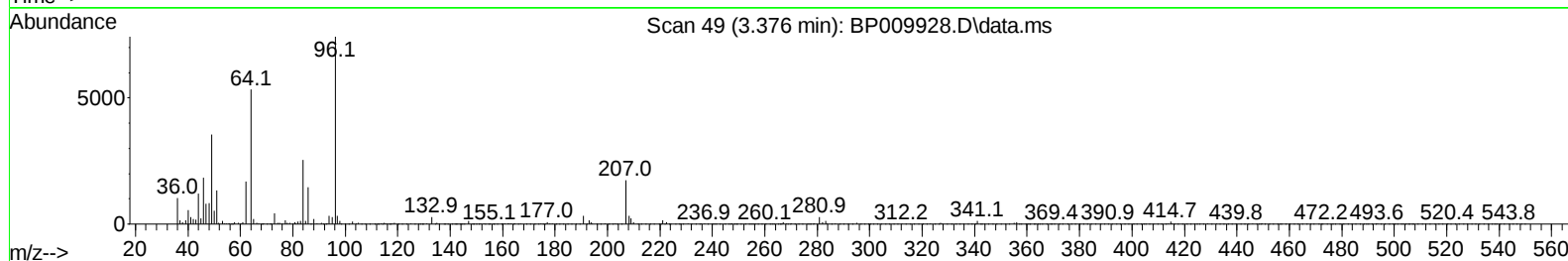
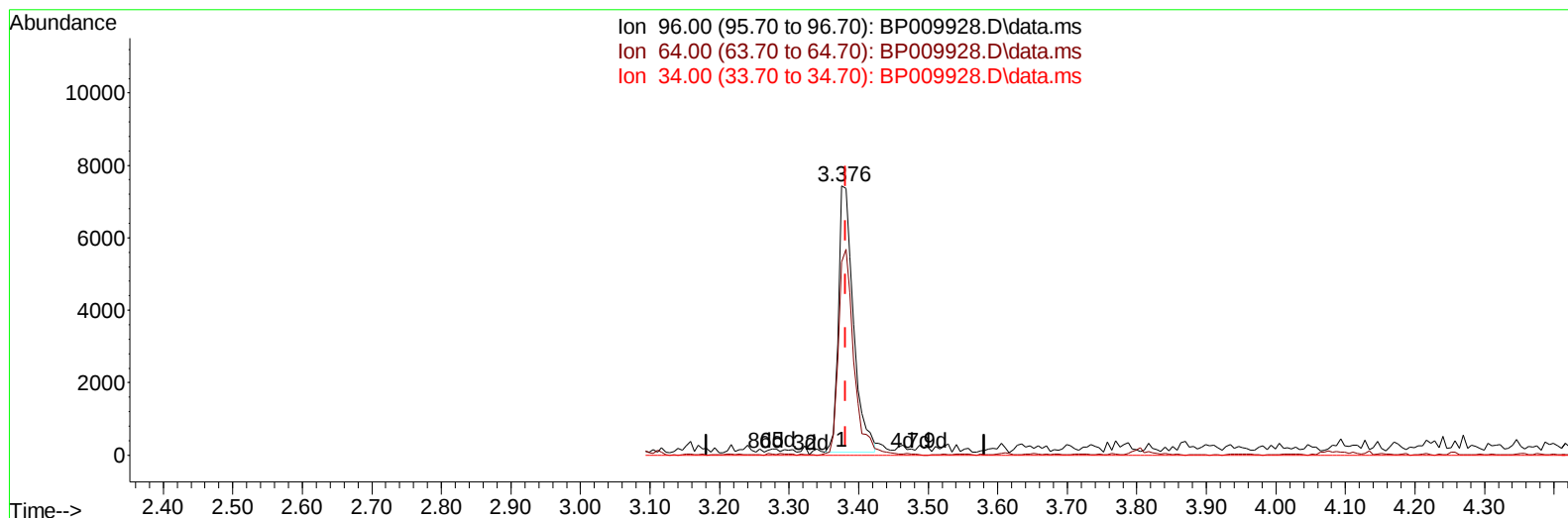
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041822\
 Data File : BP009928.D
 Acq On : 18 Apr 2022 22:18
 Operator : CG/JU
 Sample : N2421-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J76

Manual IntegrationsAPPROVED

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

Quant Time: Apr 19 00:59:21 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



TIC: BP009928.D\data.ms

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

3.376min (-0.006) 4.57 ng/uL

response 11099

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	71.86#
34.00	0.00	0.00
0.00	0.00	0.00

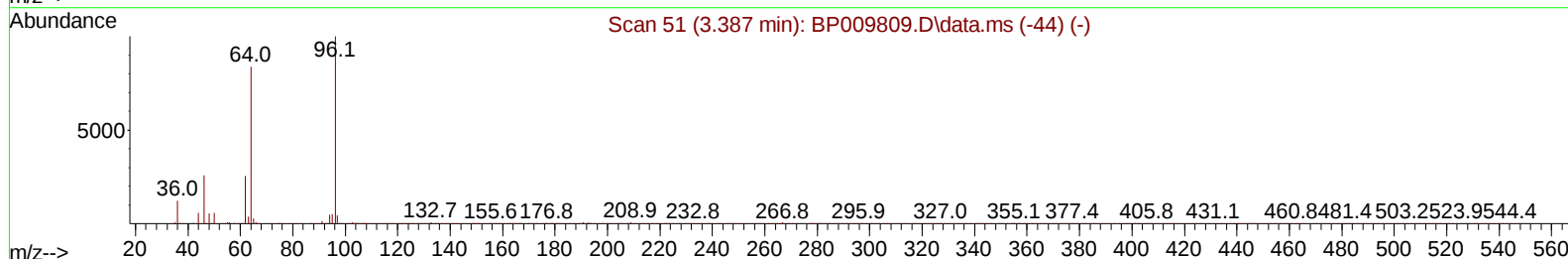
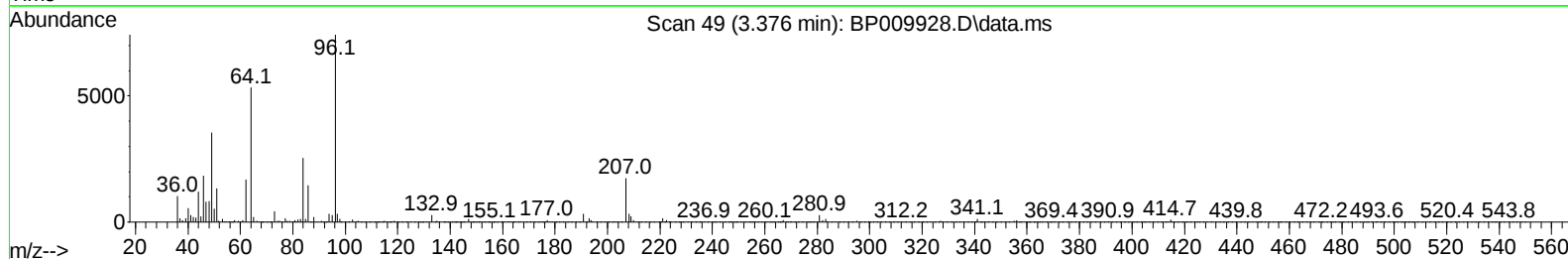
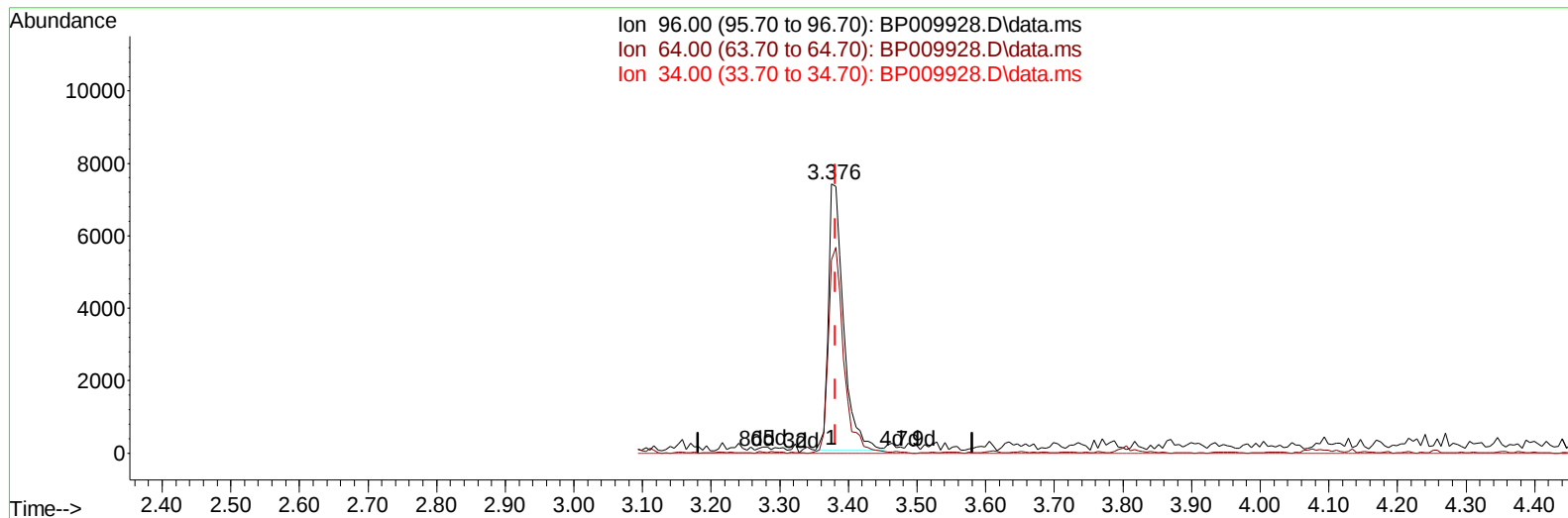
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TIC: BP009928.D\data.ms

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

3.376min (-0.006) 4.69 ng/uL m

response 11377

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	71.86#
34.00	0.00	0.00
0.00	0.00	0.00

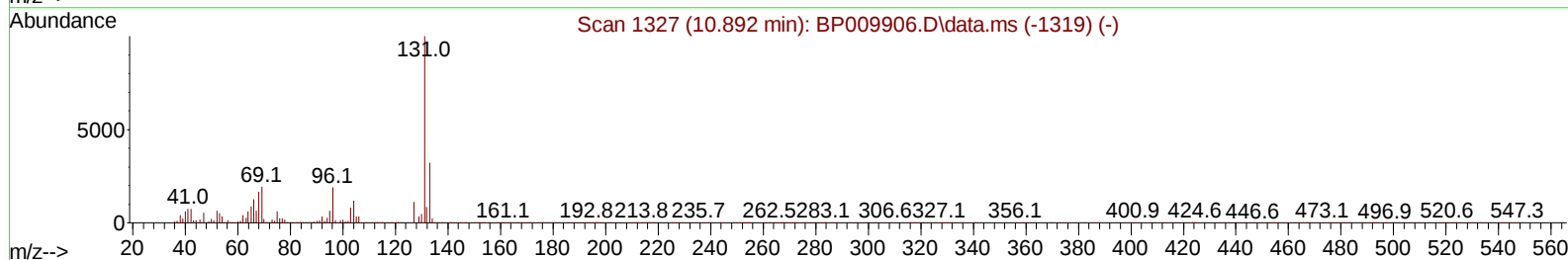
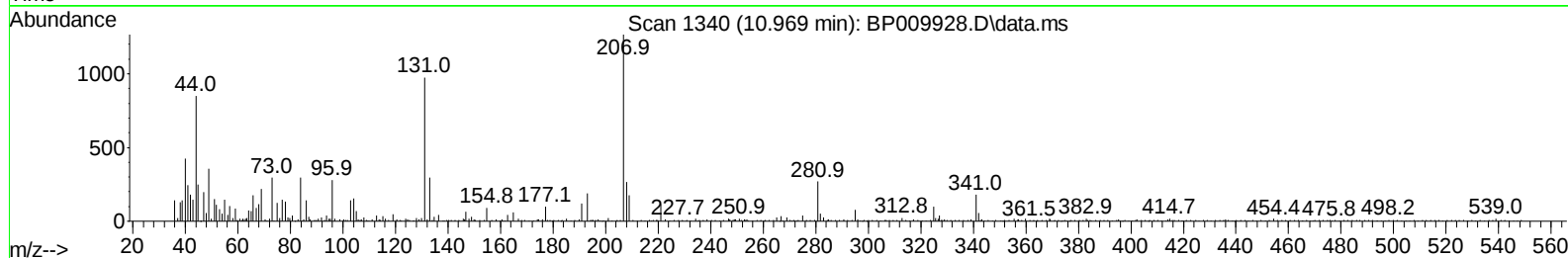
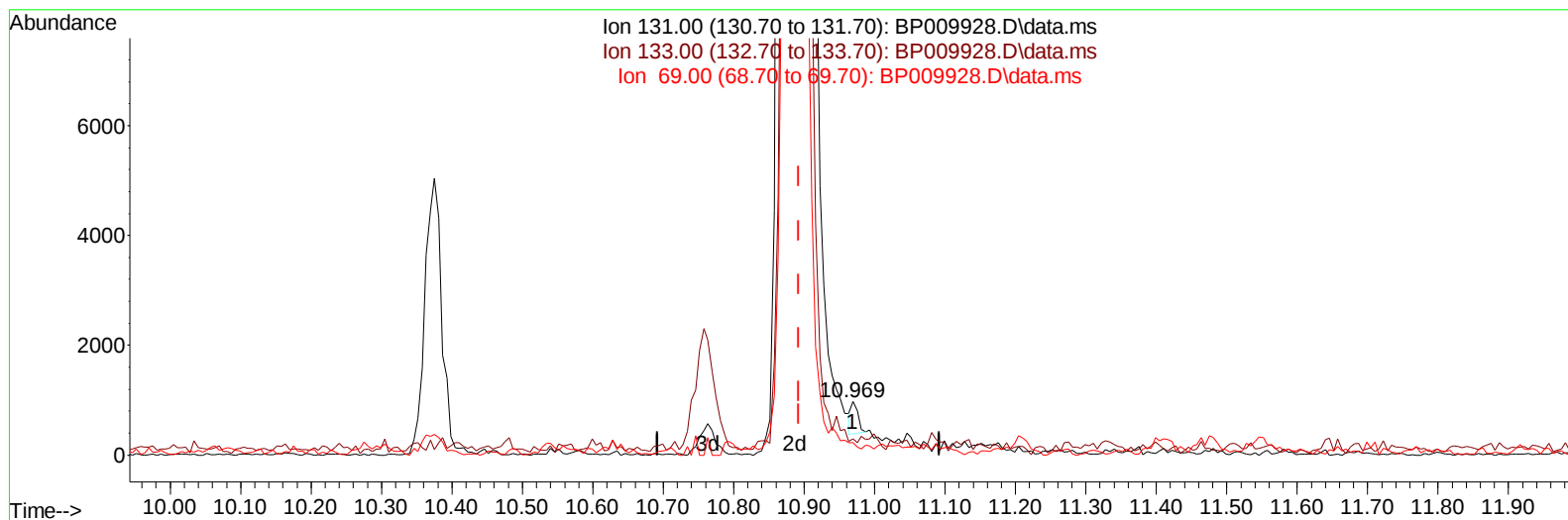
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TIC: BP009928.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.969min (+ 0.077) 0.03 ng/u1

response 377

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	30.36
69.00	26.10	22.67
0.00	0.00	0.00

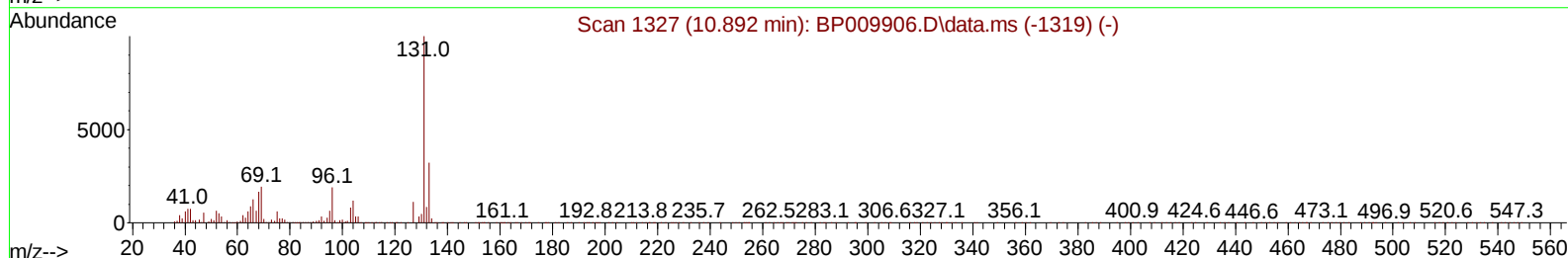
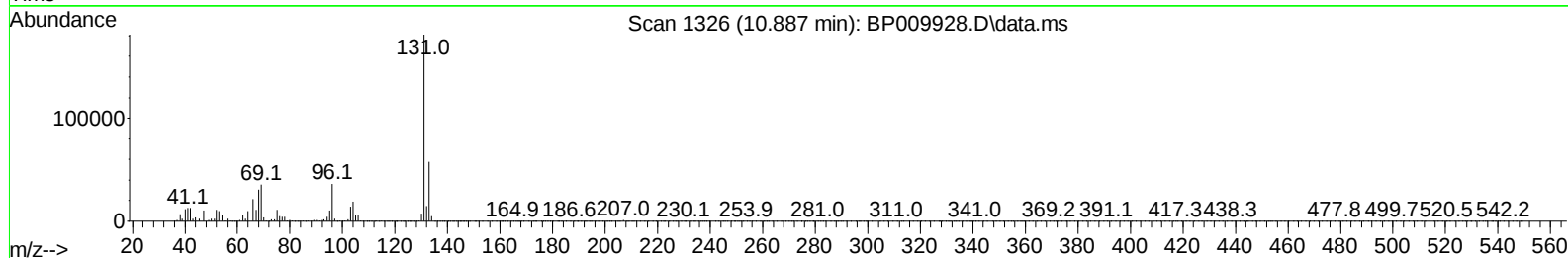
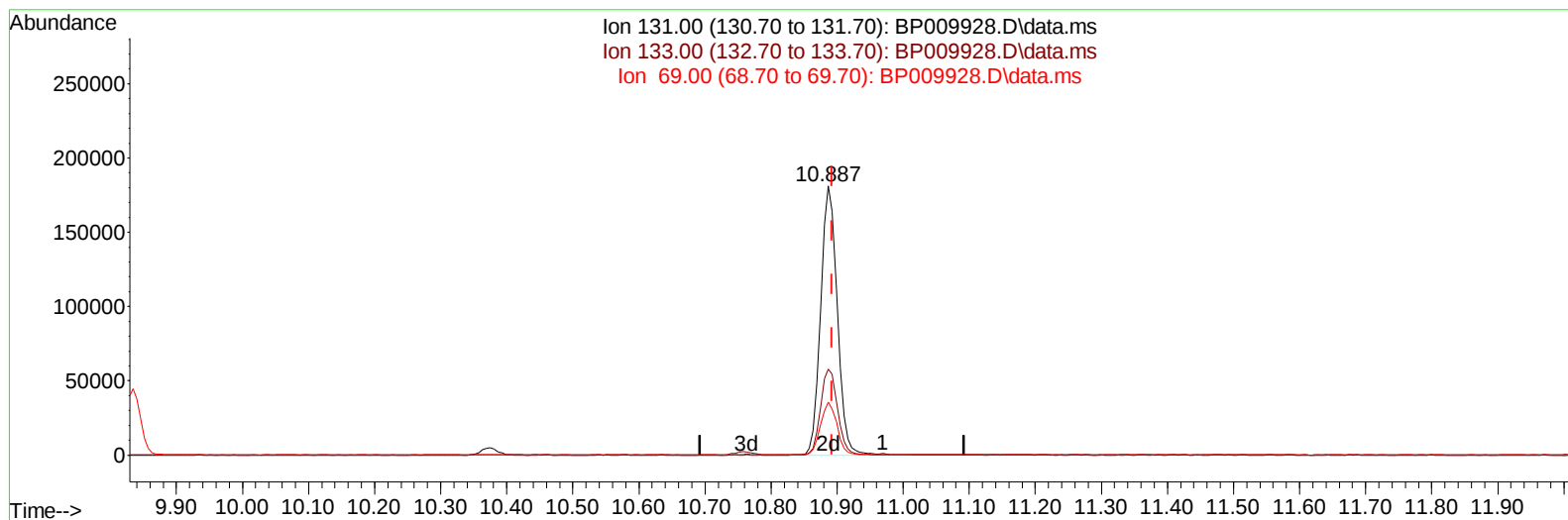
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TIC: BP009928.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.887min (-0.006) 27.30 ng/ul m

response 316321

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	31.85
69.00	26.10	19.60#
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	7.952	152	108300	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.758	136	491685	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.581	164	367905	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.334	188	831653	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.416	240	792645	20.000	ng/ul	0.00
88) Perylene-d12	23.874	264	696805	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	11377m	4.688	ng/uL	0.00
4) Pyridine-d5	3.799	84	73921	11.049	ng/ul	0.00
7) Phenol-d5	7.105	99	61182	6.435	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	184535	32.601	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	172721	24.082	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	124856	15.730	ng/ul	-0.01
21) Nitrobenzene-d5	9.111	128	125066	35.273	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	125790	32.359	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.375	165	226656	27.864	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	316321m	27.303	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	979053	33.878	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	1067575	31.181	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	29217	5.575	ng/ul	0.00
60) Fluorene-d10	15.575	176	854331	35.141	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	153114	31.078	ng/ul	0.00
73) Anthracene-d10	17.434	188	1503650	37.762	ng/ul	0.00
81) Pyrene-d10	19.663	212	1829952	42.271	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	1436732	39.593	ng/ul	-0.01

Target Compounds Qvalue

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

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