

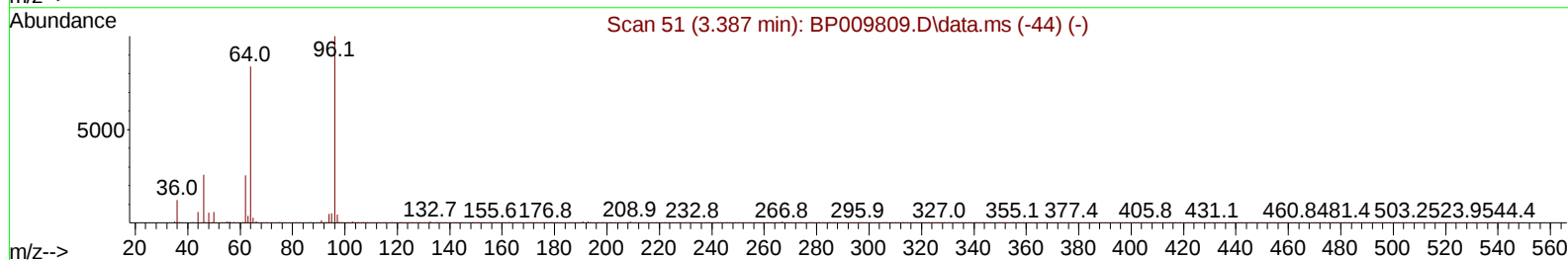
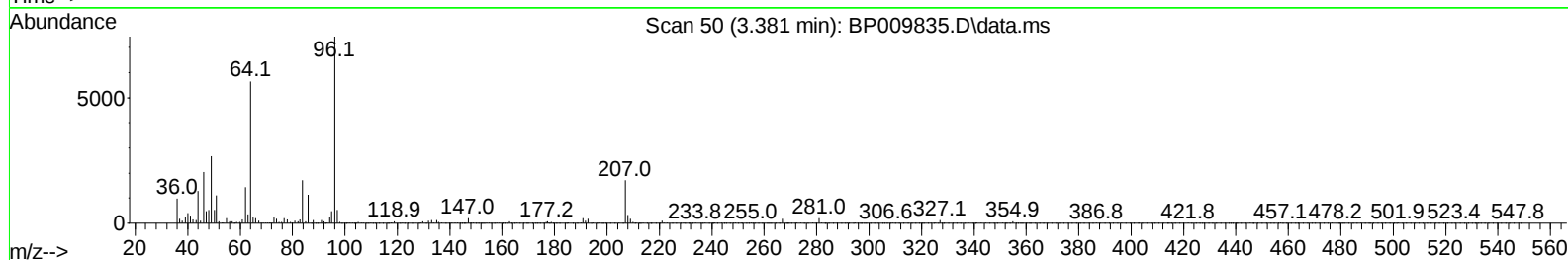
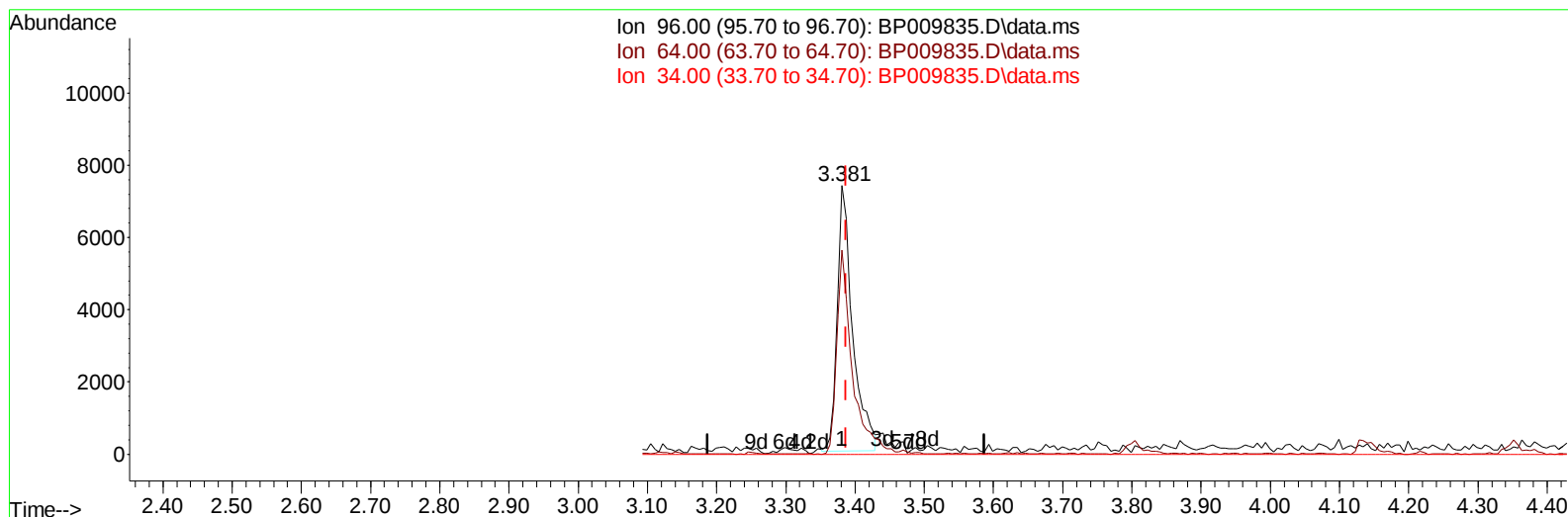
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041422\
 Data File : BP009835.D
 Acq On : 14 Apr 2022 19:13
 Operator : CG/JU
 Sample : N2293-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNP4

Manual IntegrationsAPPROVED

Quant Time: Apr 15 02:31:00 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: BP009835.D\data.ms

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

3.381min (-0.006) 5.12 ng/uL

response 11243

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	76.09#
34.00	0.00	0.00
0.00	0.00	0.00

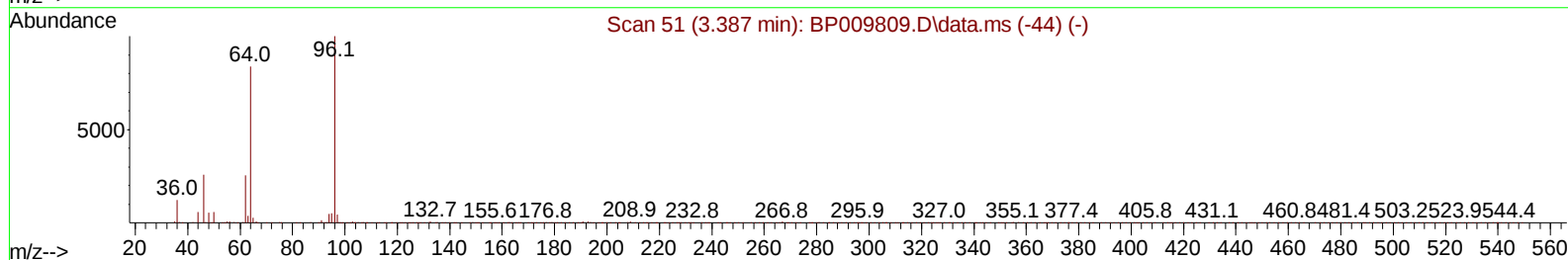
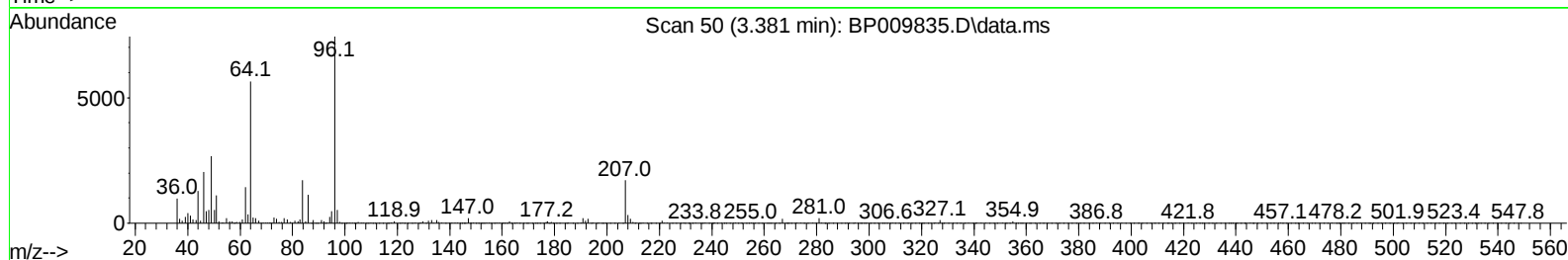
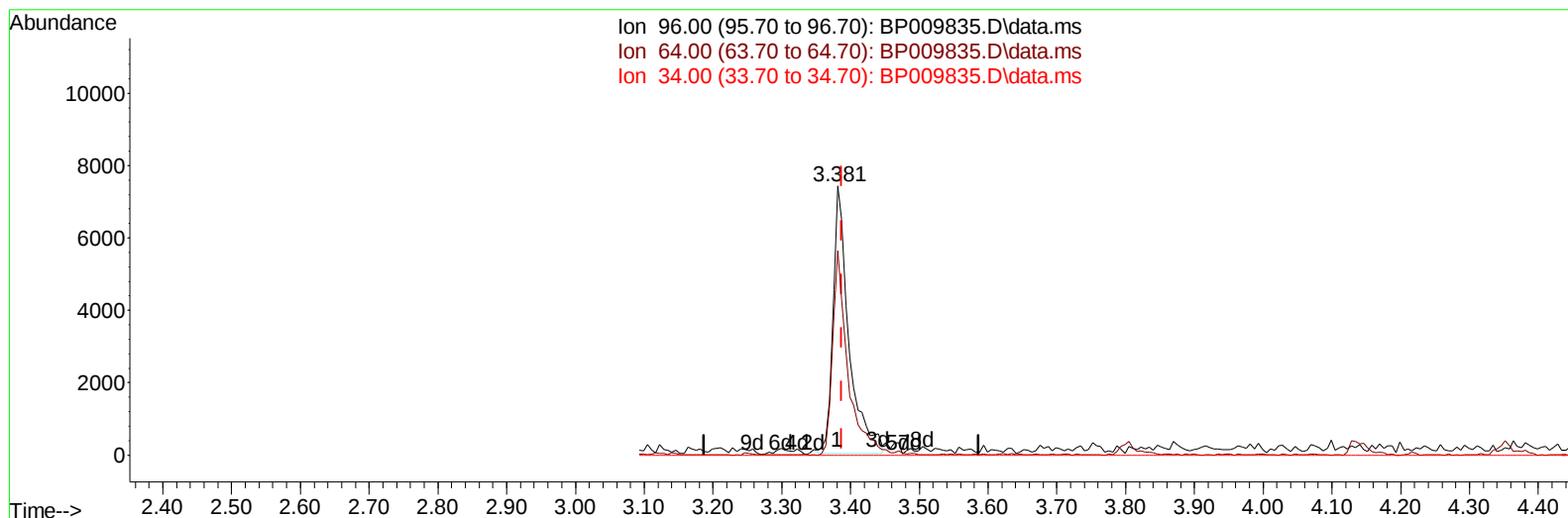
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TIC: BP009835.D\data.ms

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

3.381min (-0.006) 5.43 ng/uL m

response 11922

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	76.09#
34.00	0.00	0.00
0.00	0.00	0.00

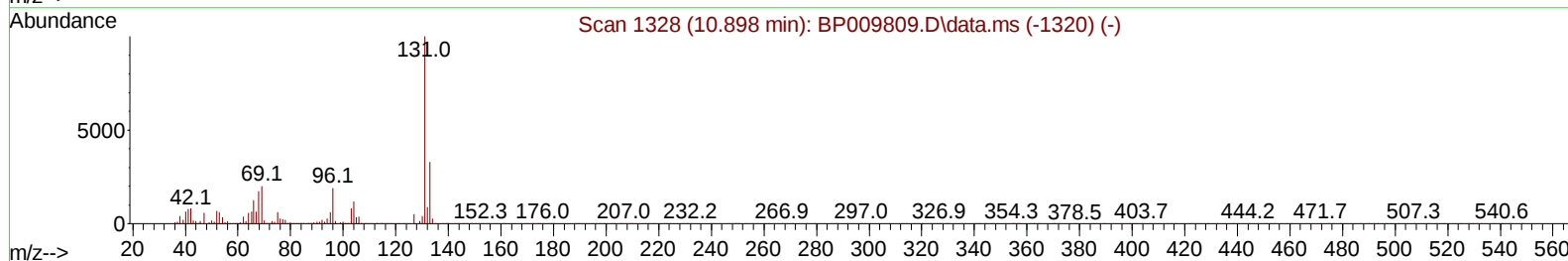
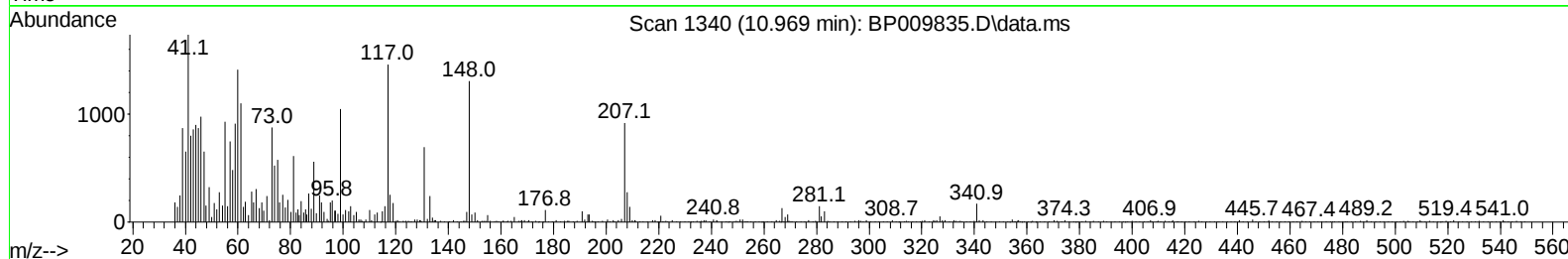
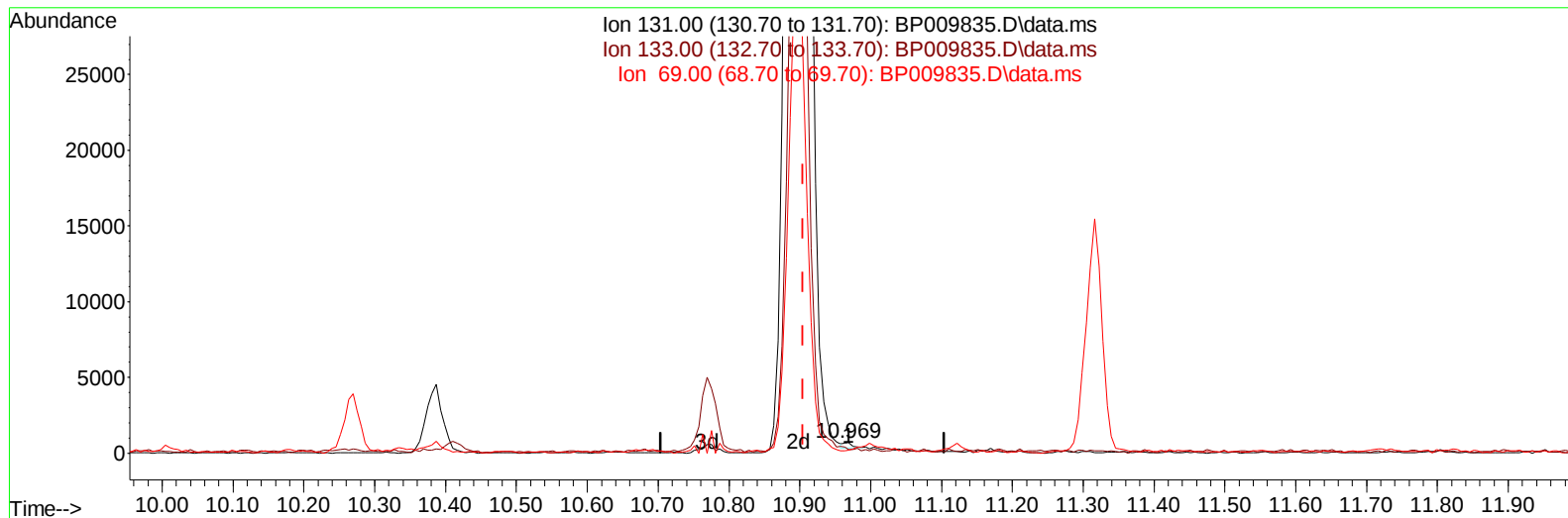
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TIC: BP009835.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.969min (+ 0.065) 0.03 ng/u1

response 308

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	34.83
69.00	26.10	26.16
0.00	0.00	0.00

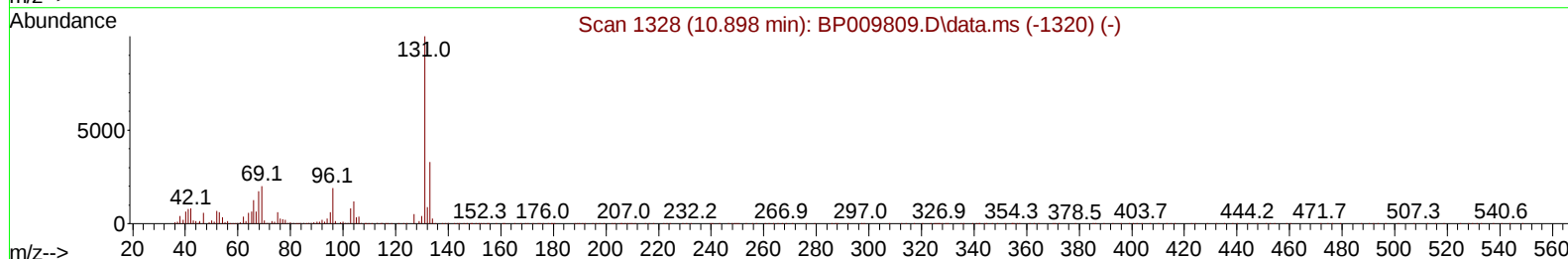
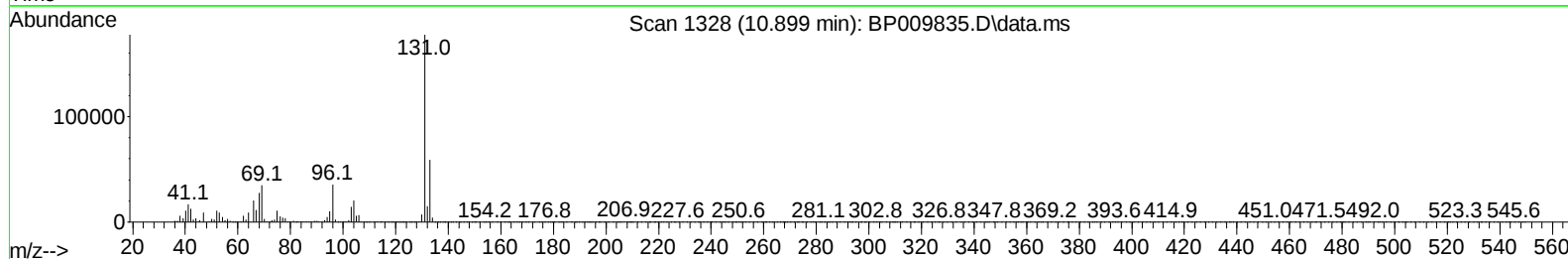
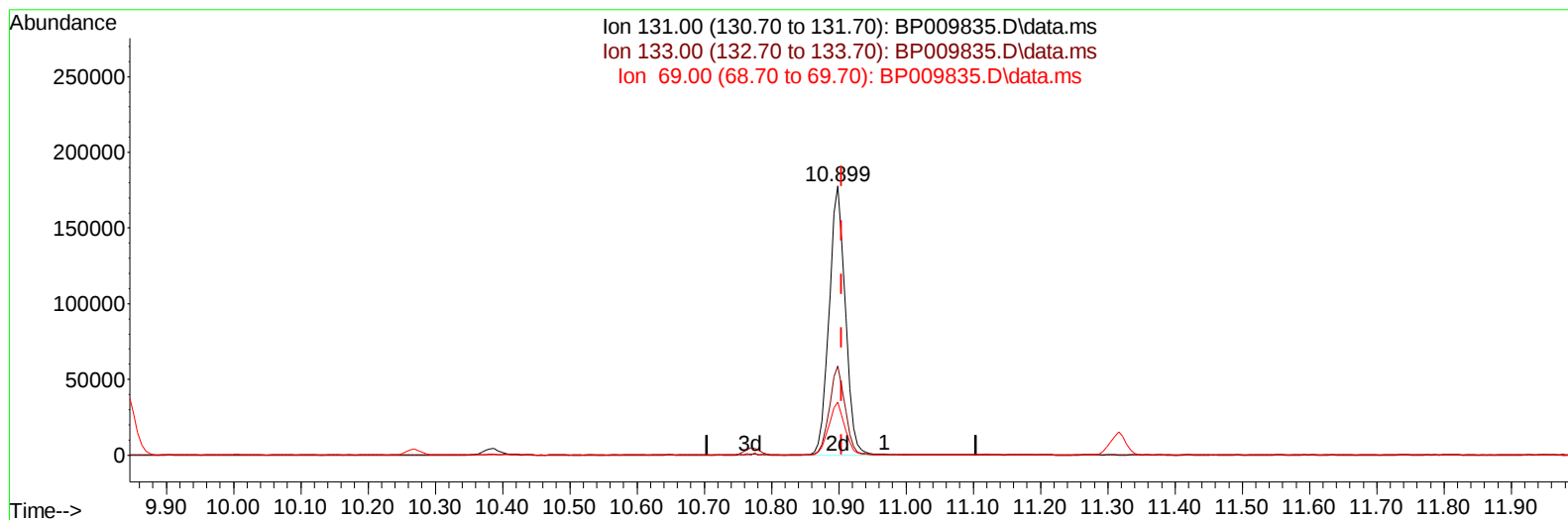
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TIC: BP009835.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.899min (-0.006) 28.91 ng/ul m

response 298100

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	33.10
69.00	26.10	19.73#
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	97977	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	437585	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	314772	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	729812	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.422	240	836737	20.000	ng/ul	#-0.01
88) Perylene-d12	23.880	264	867723	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	11922m	5.431	ng/uL	0.00
4) Pyridine-d5	3.799	84	157179	25.968	ng/ul	0.00
7) Phenol-d5	7.116	99	237821	27.648	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	154194	30.111	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	186223	28.701	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	212571	29.603	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	99903	31.659	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.846	143	107963	31.207	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	205705	28.415	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.899	131	298100m	28.911	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	757143	30.622	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	840125	28.680	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.769	143	124304	27.721	ng/ul	-0.01
60) Fluorene-d10	15.581	176	653875	31.436	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.692	200	100796	23.314	ng/ul	0.00
73) Anthracene-d10	17.439	188	1053946	30.162	ng/ul	-0.01
81) Pyrene-d10	19.669	212	1329068	29.083	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1384485	30.638	ng/ul	-0.01
Target Compounds						
29) 2,4-Dichlorophenol	10.410	162	22329	3.181	ng/ul#	Qvalue 77

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

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