

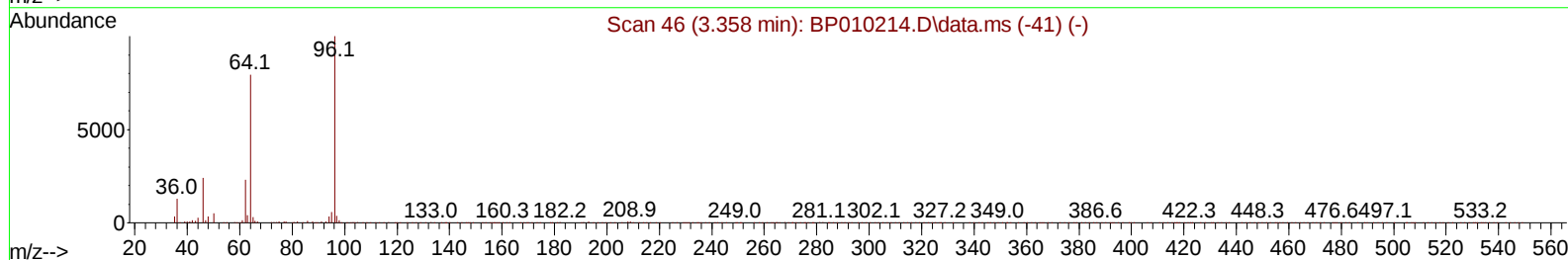
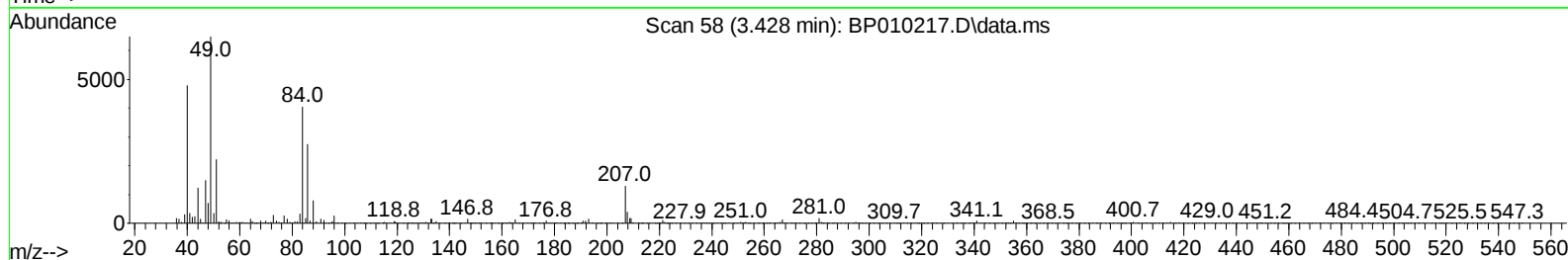
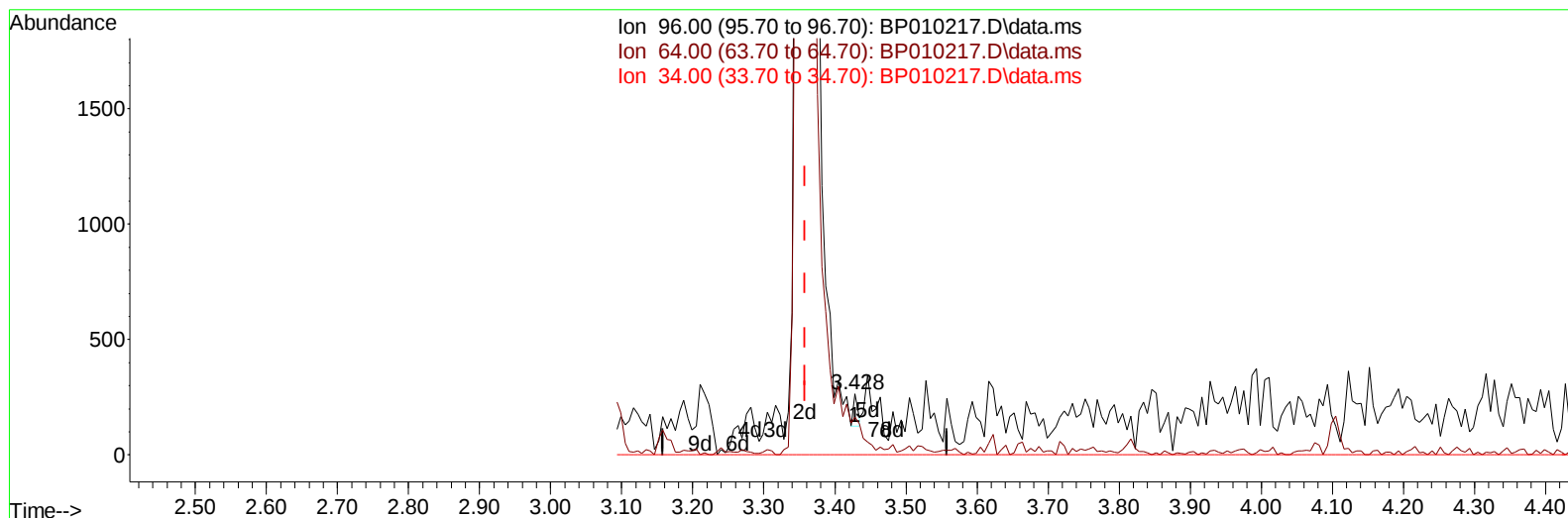
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
 Data File : BP010217.D
 Acq On : 03 May 2022 21:12
 Operator : CG/JU
 Sample : N2670-02
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COW01

Manual IntegrationsAPPROVED

Quant Time: May 04 01:25:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 00:58:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/04/2022
 Supervised By :Yogesh Patel 05/05/2022



TIC: BP010217.D\data.ms

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

3.428min (+ 0.071) 0.02 ng/uL

response 62

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	62.12
34.00	0.00	0.00
0.00	0.00	0.00

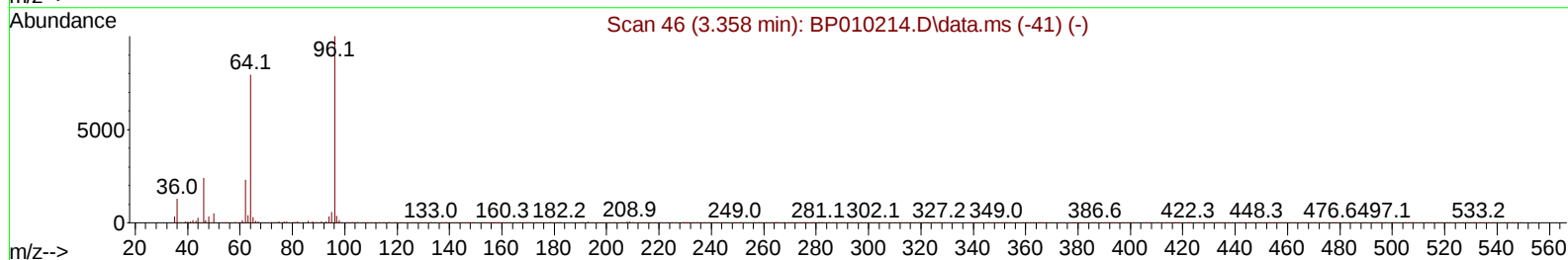
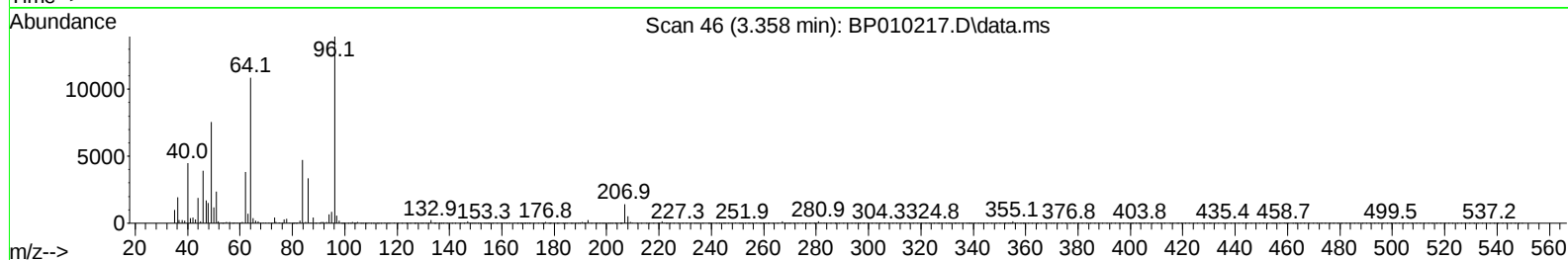
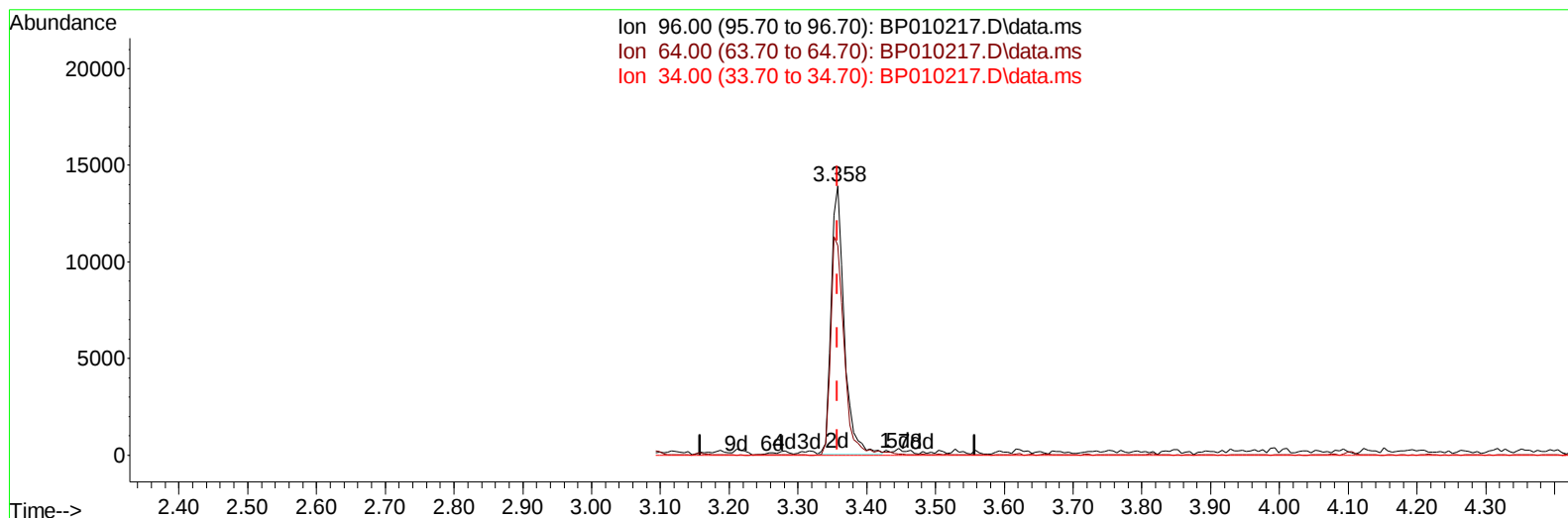
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TIC: BP010217.D\data.ms

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

3.358min (-0.000) 5.44 ng/uL m

response 18344

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.02#
34.00	0.00	0.00
0.00	0.00	0.00

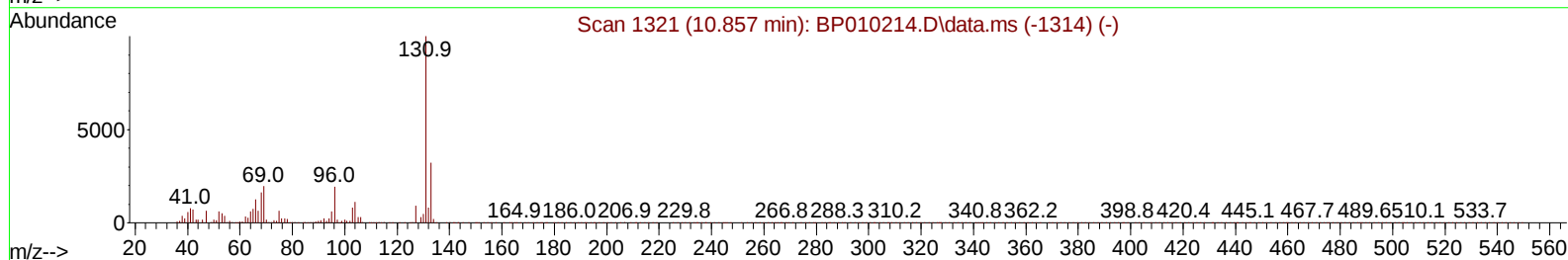
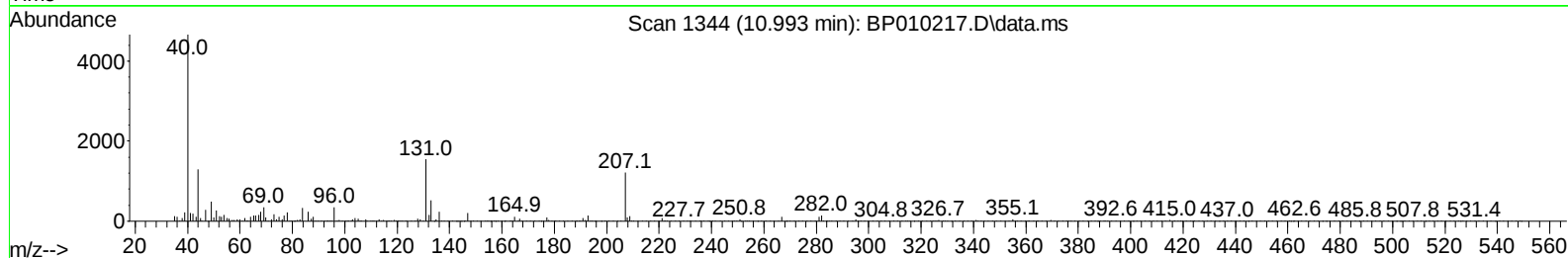
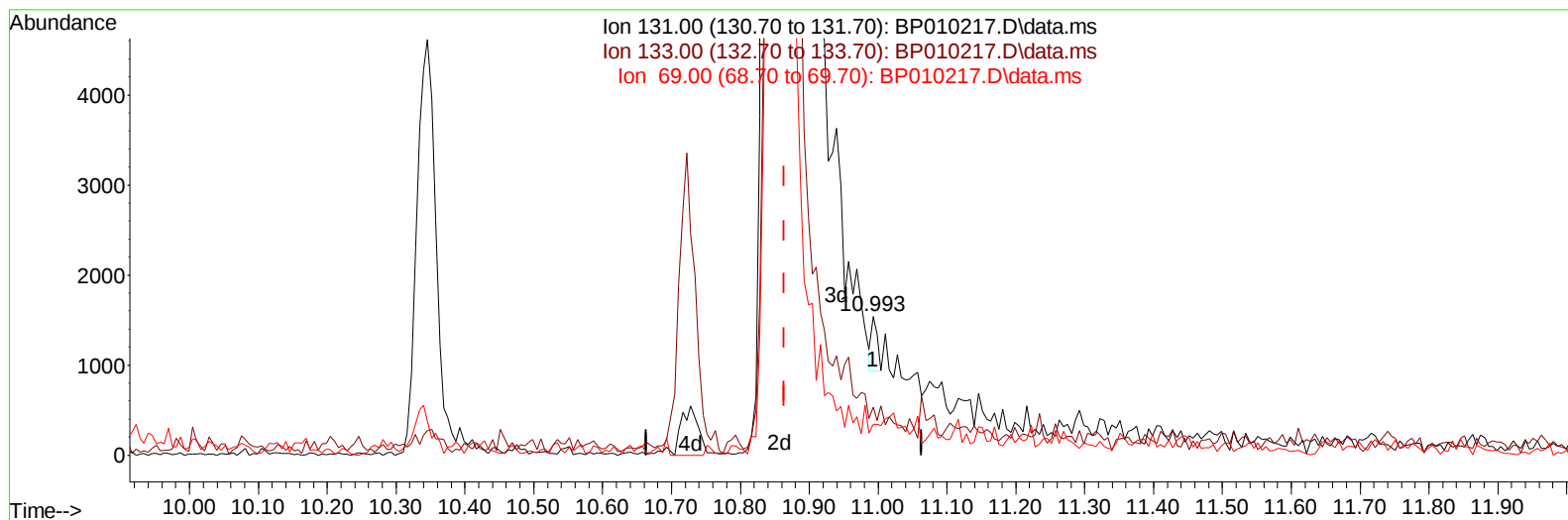
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TIC: BP010217.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.993min (+ 0.129) 0.02 ng/u1

response 359

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	34.33
69.00	26.10	22.19
0.00	0.00	0.00

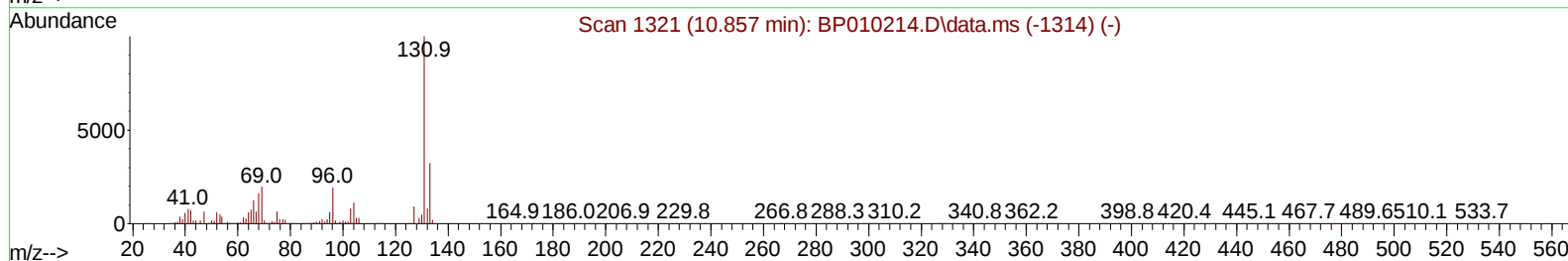
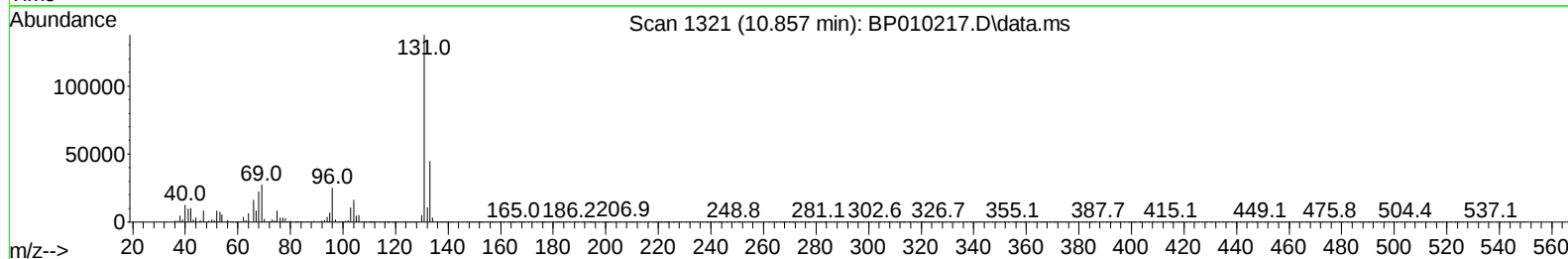
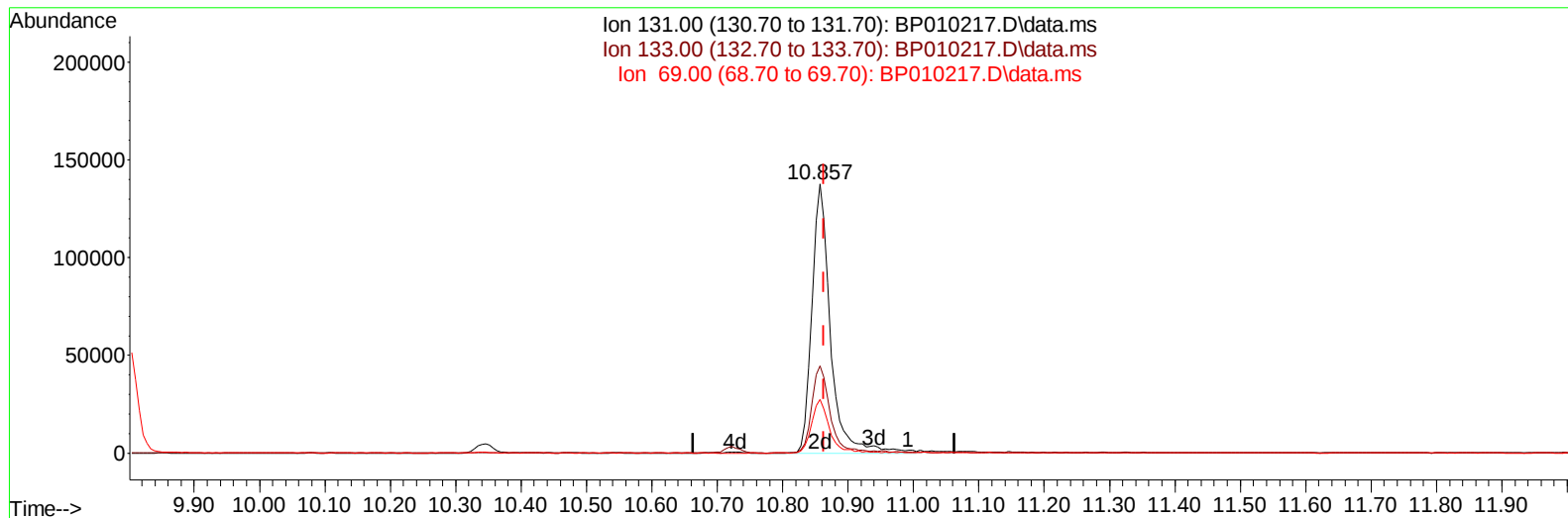
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Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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TIC: BP010217.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.857min (-0.006) 18.54 ng/ul m

response 271714

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.47
69.00	26.10	19.88#
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	7.916	152	147108	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.722	136	604376	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	367329	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	721336	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	649150	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	615232	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	18344m	5.438	ng/uL	0.00
4) Pyridine-d5	3.787	84	28449	2.998	ng/ul	0.01
7) Phenol-d5	7.081	99	66018	4.934	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	263184	32.779	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	224663	22.671	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	135103	12.186	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	151941	31.609	ng/ul	0.00
24) 2-Nitrophenol-d4	9.805	143	158409	30.175	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	245309	24.003	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	271714m	18.542	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	912389	31.931	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1034044	29.901	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	15785	3.001	ng/ul	0.01
60) Fluorene-d10	15.545	176	779425	31.710	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	110714	24.273	ng/ul	0.00
73) Anthracene-d10	17.398	188	1251090	35.850	ng/ul	-0.01
81) Pyrene-d10	19.633	212	1422523	38.533	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1244948	38.598	ng/ul	-0.01

Target Compounds Qvalue

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

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