

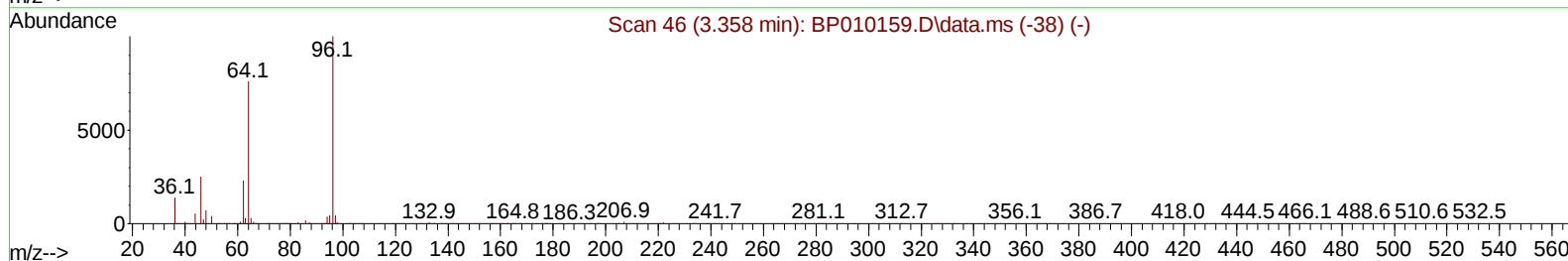
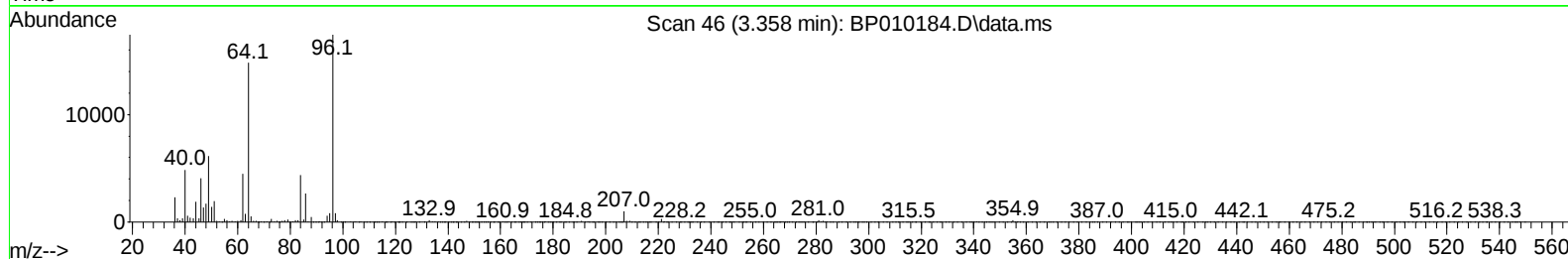
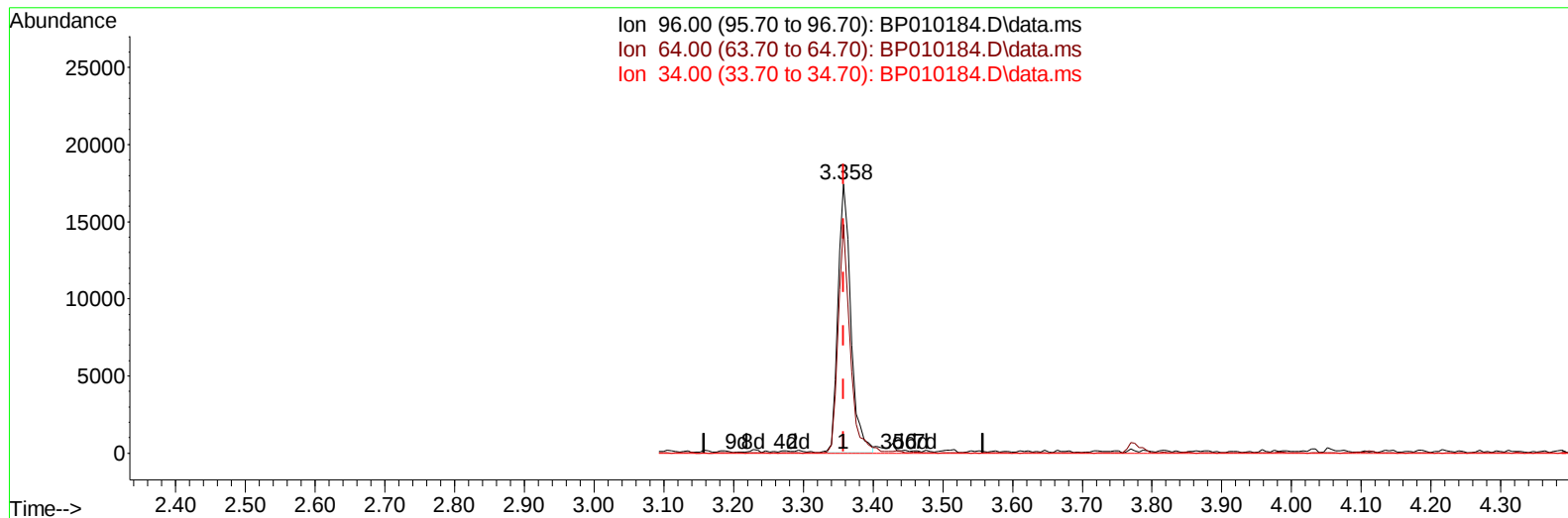
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050222\
Data File : BP010184.D
Acq On : 02 May 2022 22:02
Operator : CG/JU
Sample : PB144527BL
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK527

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/03/2022
Supervised By :Yogesh Patel 05/05/2022

Quant Time: May 03 01:38:42 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



TIC: BP010184.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.358min (+ 0.000) 6.23 ng/uL****response 22156**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.98#
34.00	0.00	0.00
0.00	0.00	0.00

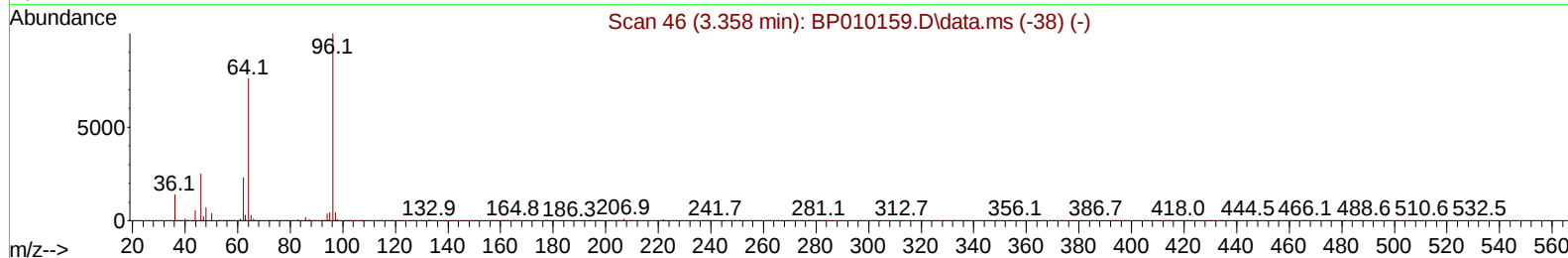
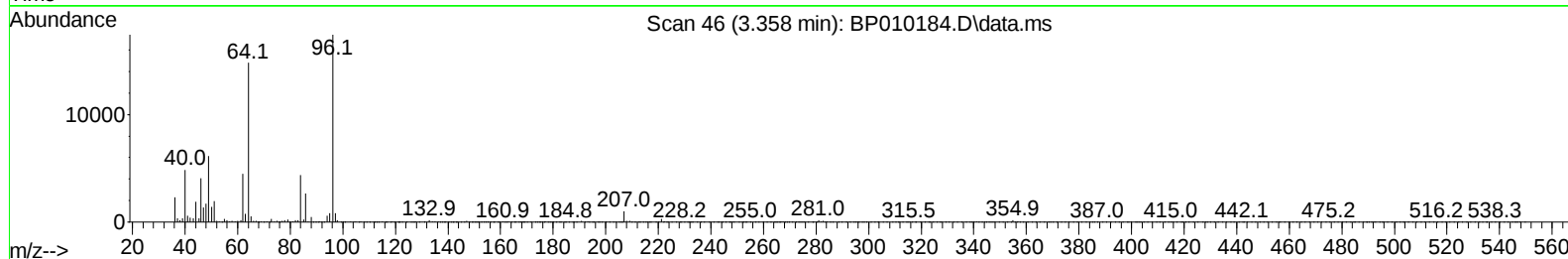
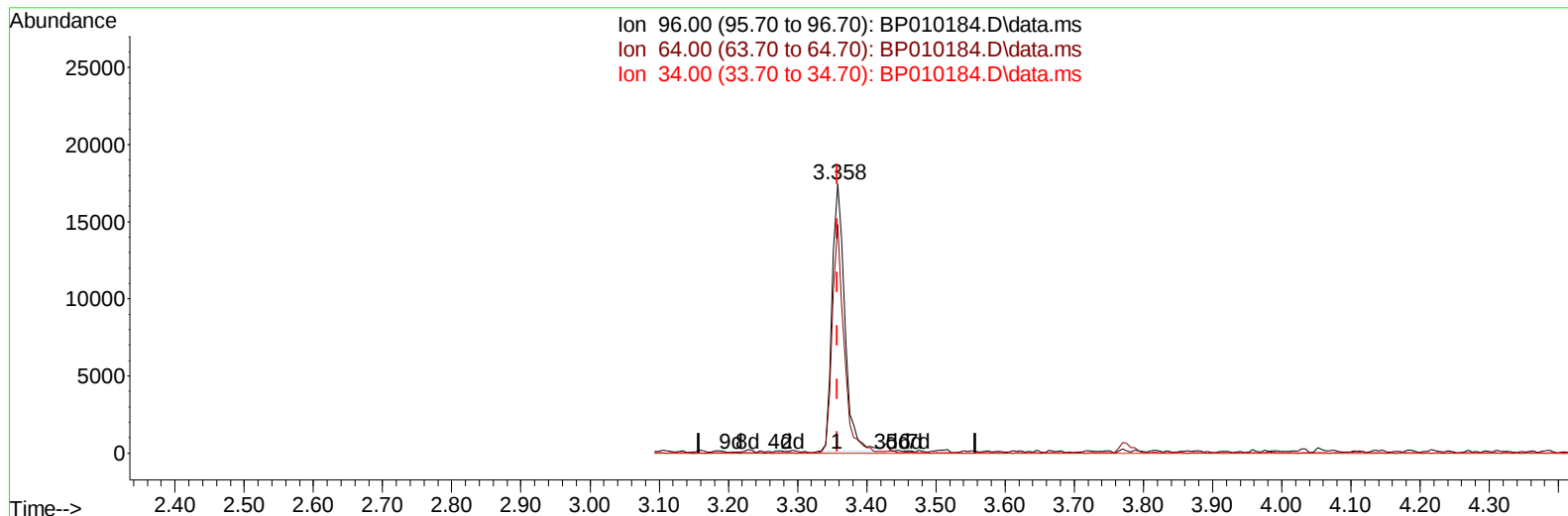
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TIC: BP010184.D\data.ms

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

3.358min (+ 0.000) 6.24 ng/uL m

response 22191

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.98#
34.00	0.00	0.00
0.00	0.00	0.00

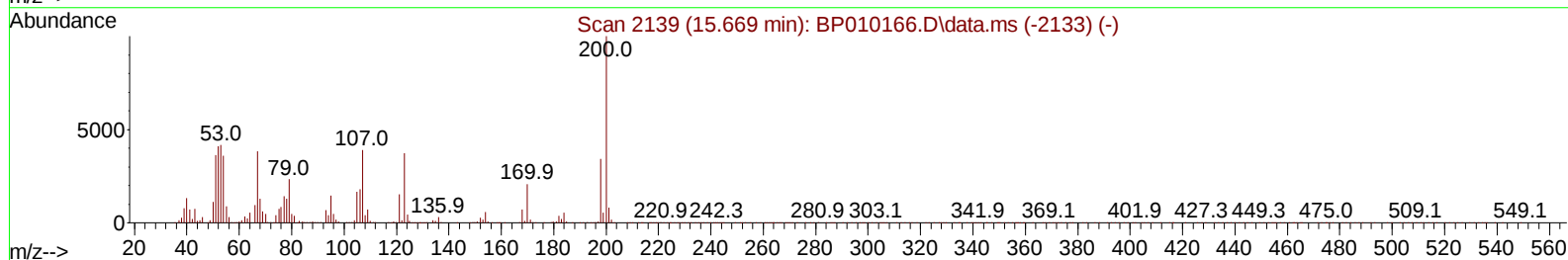
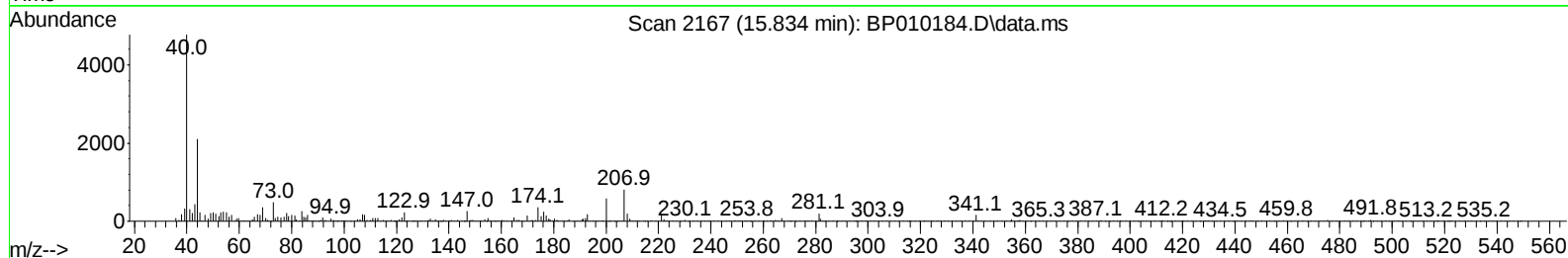
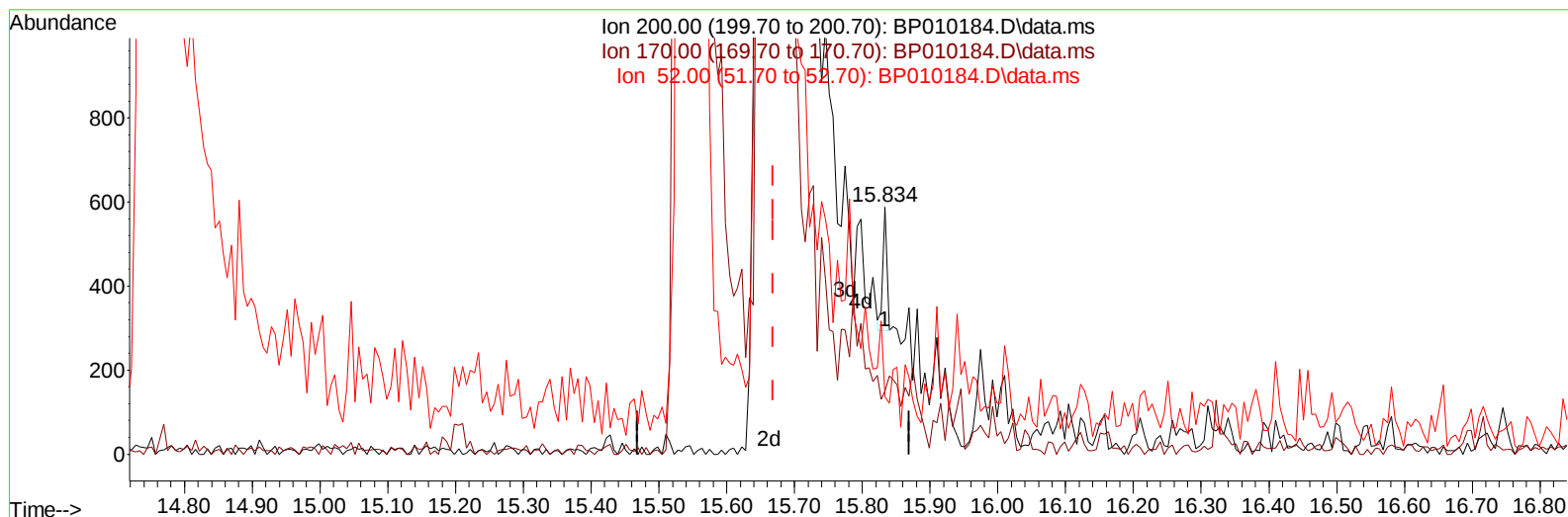
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TIC: BP010184.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.834min (+ 0.165) 0.03 ng/u1****response 120**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	25.51
52.00	21.20	23.13
0.00	0.00	0.00

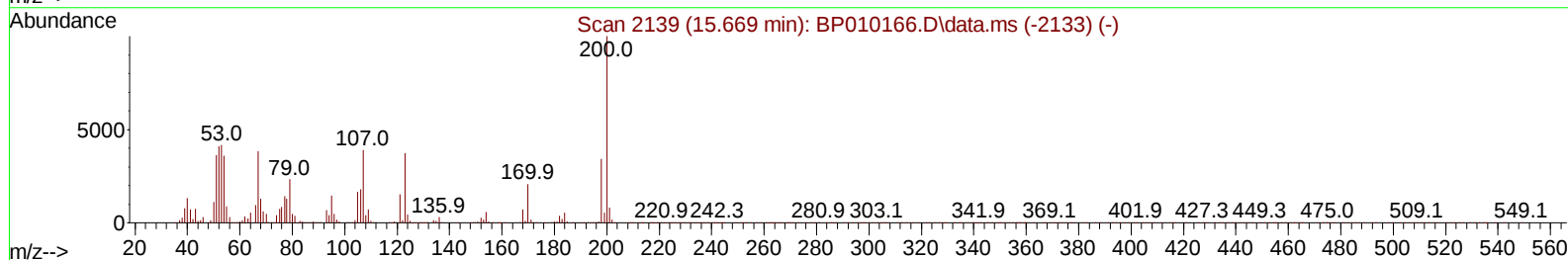
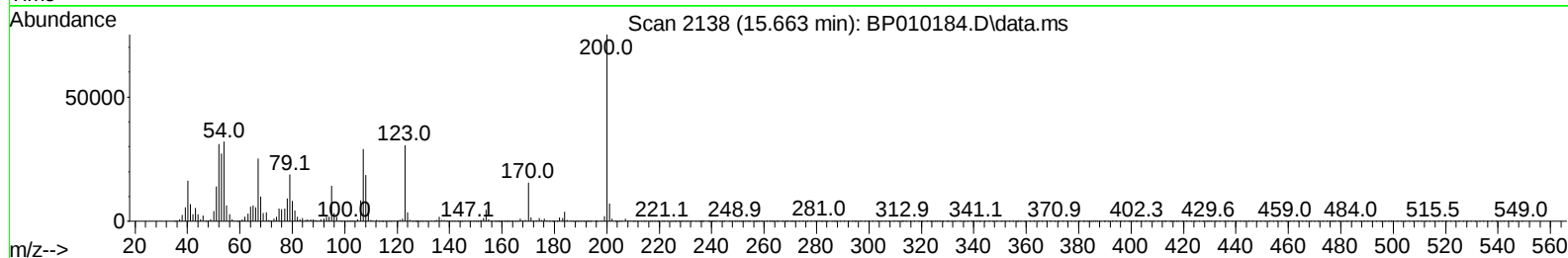
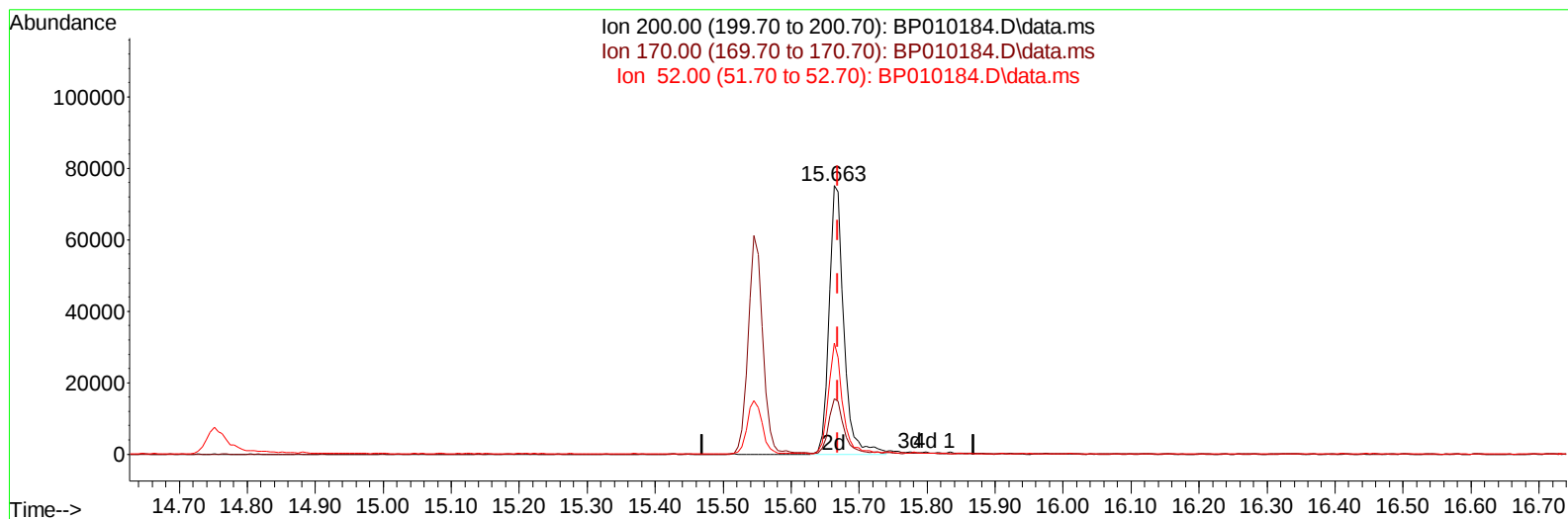
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BP010184.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.663min (-0.006) 24.25 ng/ul m

response 112529

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	20.75
52.00	21.20	41.28#
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	7.922	152	155114	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	644089	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	391378	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	733985	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	619837	20.000	ng/ul	# 0.00
88) Perylene-d12	23.833	264	584684	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	22191m	6.238	ng/uL	0.00
4) Pyridine-d5	3.775	84	306943	30.673	ng/ul	0.00
7) Phenol-d5	7.087	99	378198	26.805	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	271299	32.046	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	319197	30.548	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	322111	27.555	ng/ul	0.00
21) Nitrobenzene-d5	9.081	128	164977	32.205	ng/ul	0.00
24) 2-Nitrophenol-d4	9.805	143	181543	32.449	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	295130	27.097	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	432093	27.669	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	953257	31.311	ng/ul	0.00
49) Acenaphthylene-d8	14.246	160	1095258	29.725	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	113303	20.215	ng/ul	0.00
60) Fluorene-d10	15.545	176	793927	30.315	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	112529m	24.246	ng/ul	0.00
73) Anthracene-d10	17.404	188	1136276	31.999	ng/ul	0.00
81) Pyrene-d10	19.633	212	1269102	36.003	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1050054	34.257	ng/ul	0.00

Target Compounds Qvalue

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

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