

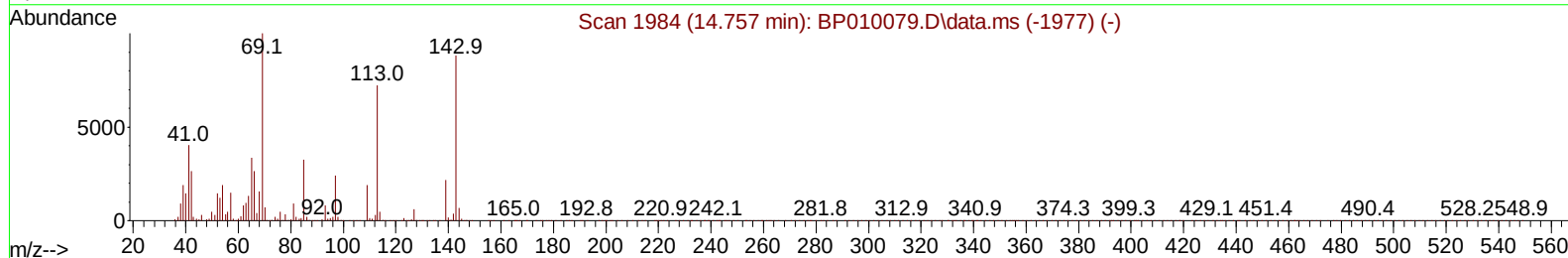
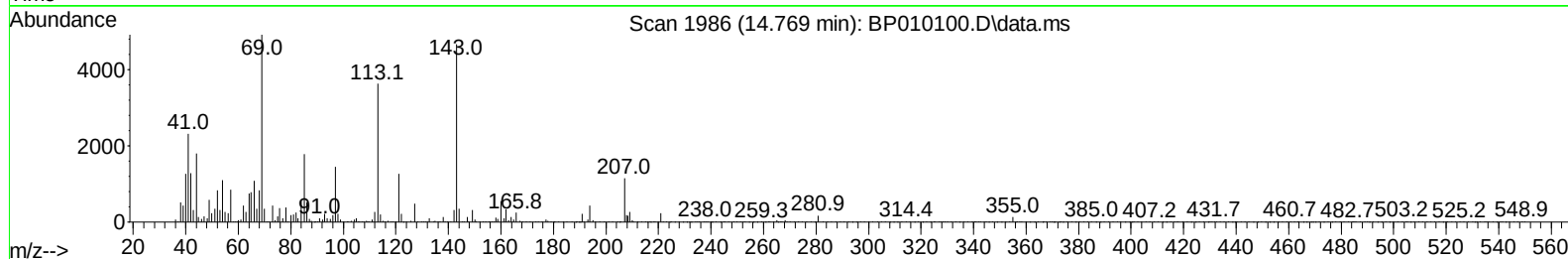
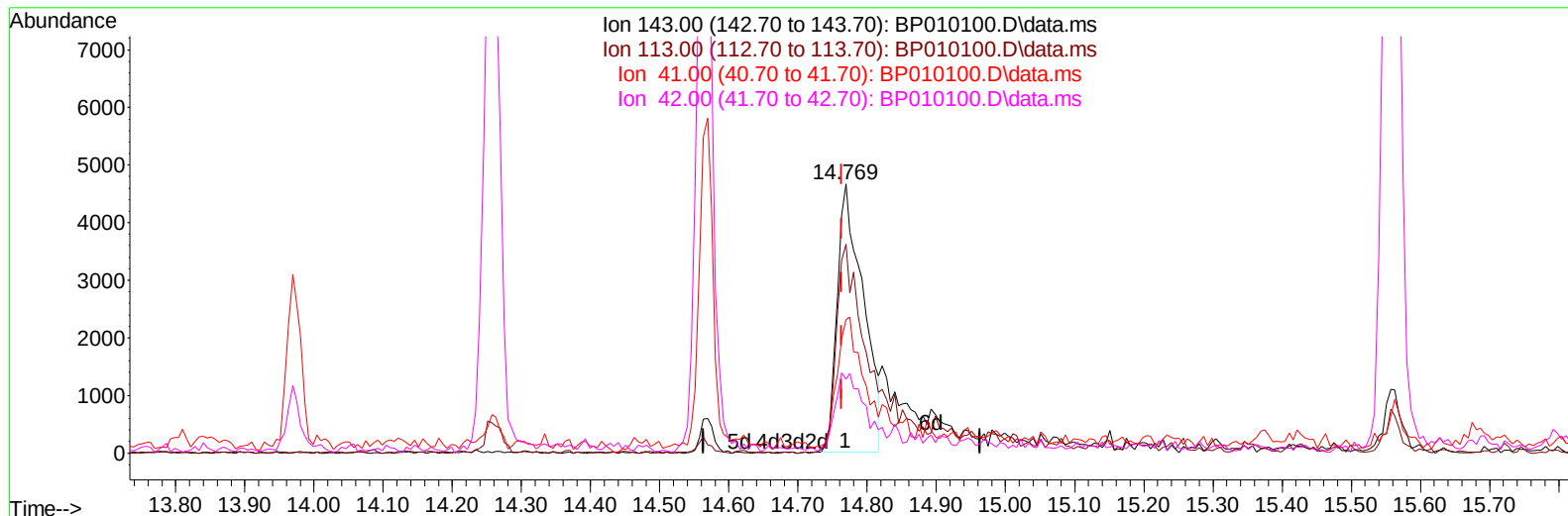
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010100.D
Acq On : 26 Apr 2022 00:57
Operator : CG/JU
Sample : N2475-15
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0BQ2

Manual IntegrationsAPPROVED

Quant Time: Apr 26 02:51:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/26/2022
Supervised By :Yogesh Patel 04/28/2022



TIC: BP010100.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 1.57 ng/ul

response 12129

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.77#
41.00	23.20	49.65#
42.00	18.00	27.54#

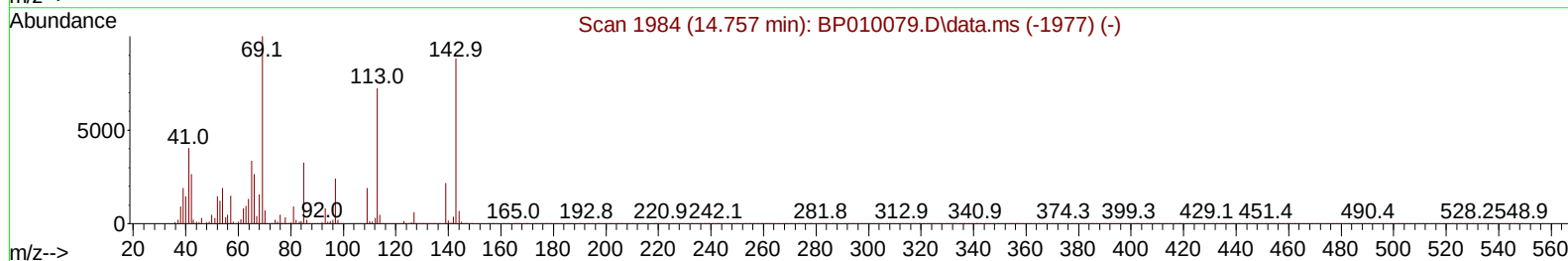
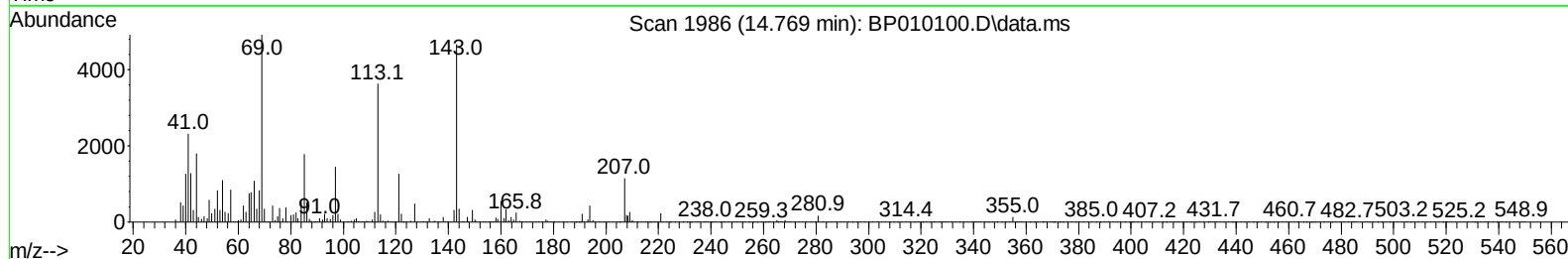
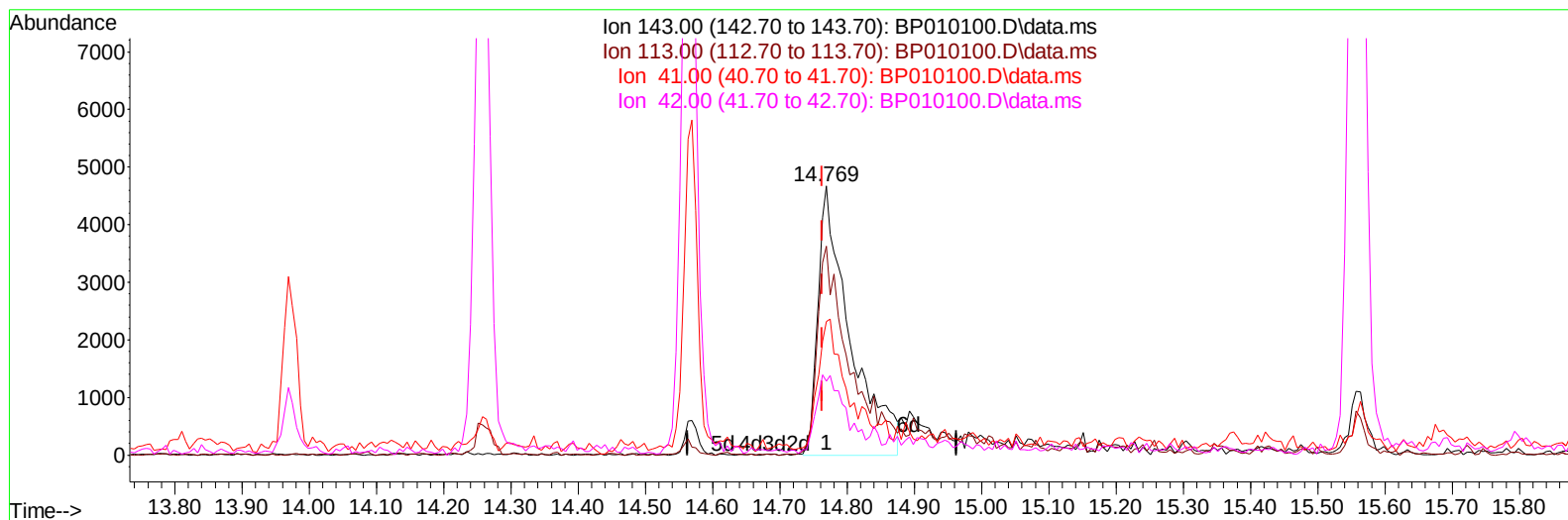
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TIC: BP010100.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 2.01 ng/ul m

response 15500

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	77.77#
41.00	23.20	49.65#
42.00	18.00	27.54#

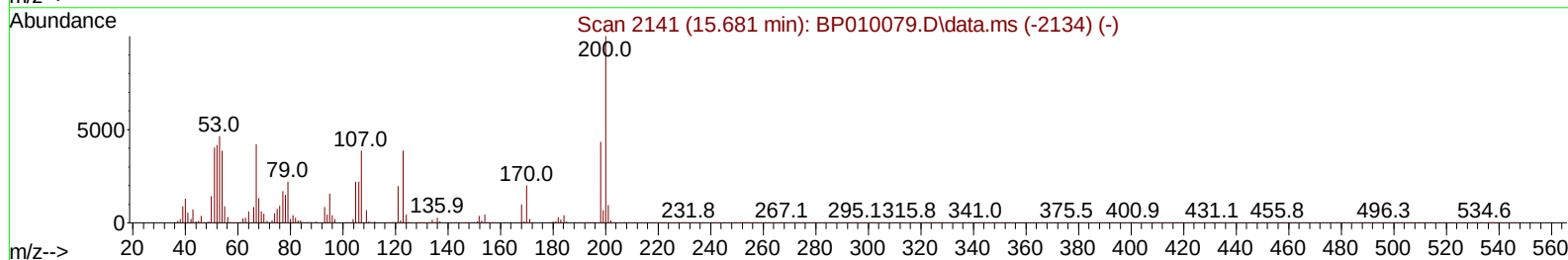
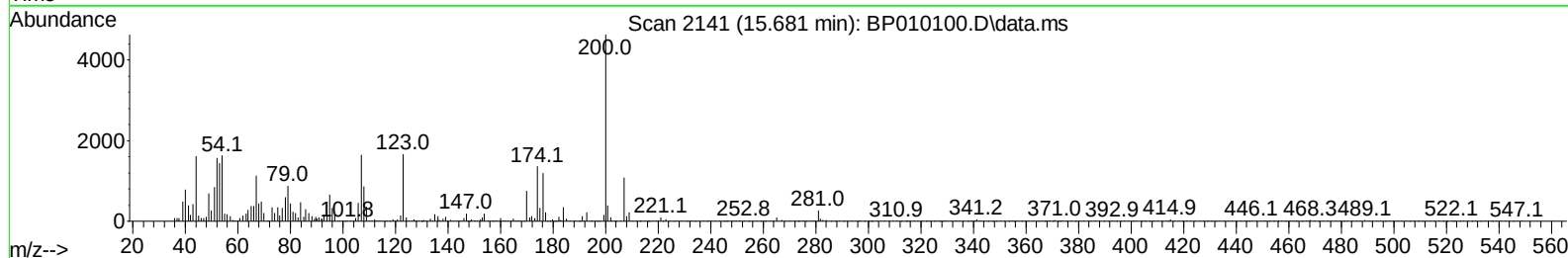
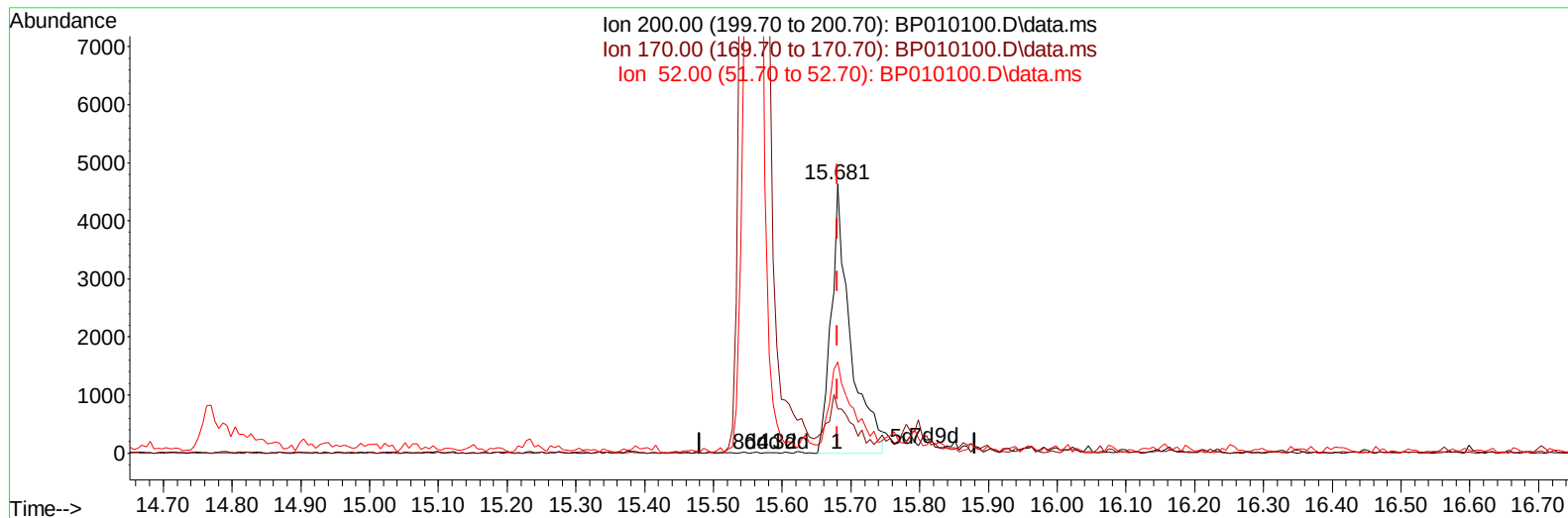
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TIC: BP010100.D\data.ms

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

15.681min (+ 0.000) 1.24 ng/ul

response 9011

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	16.32#
52.00	21.20	33.86#
0.00	0.00	0.00

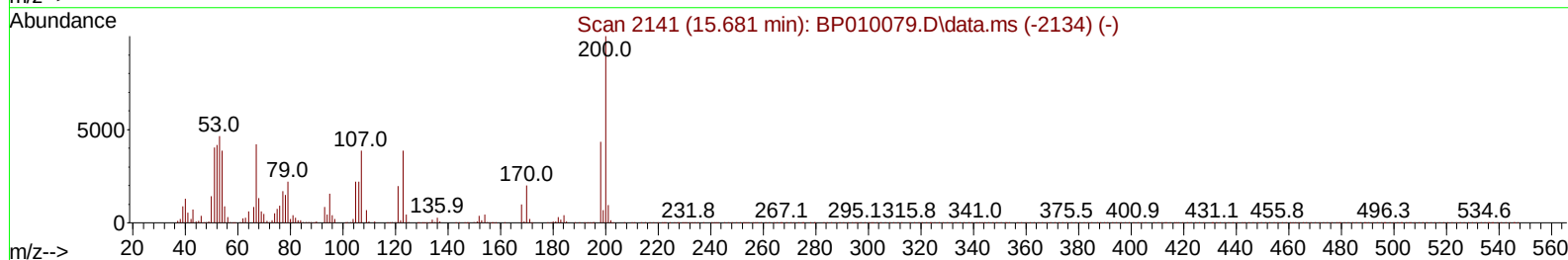
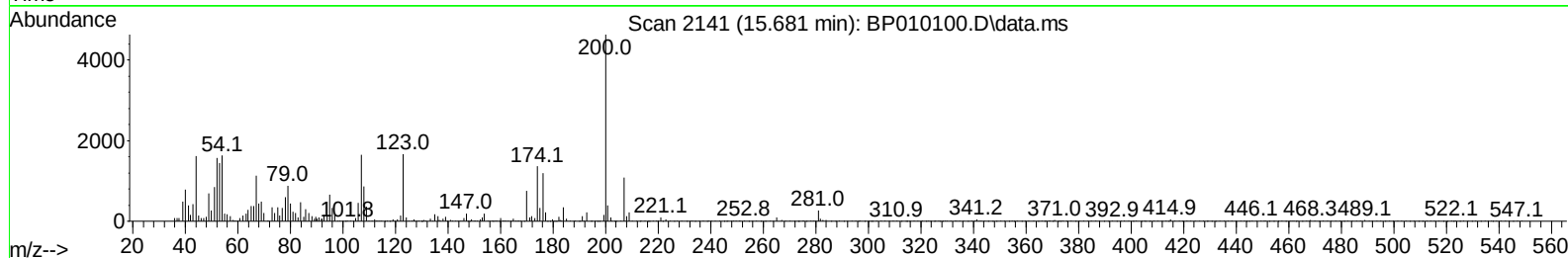
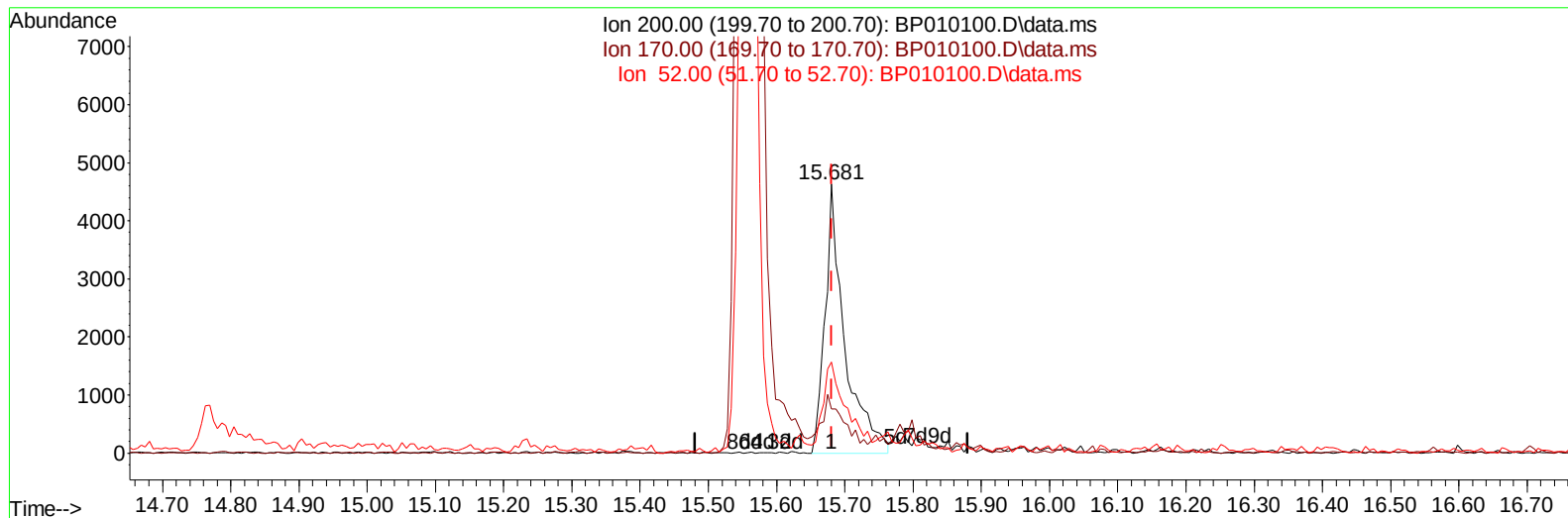
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TIC: BP010100.D\data.ms

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

15.681min (+ 0.000) 1.27 ng/ul m

response 9275

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.00	16.32#
52.00	21.20	33.86#
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	7.934	152	182764	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	786871	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	539616	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	1152667	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	958645	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	702654	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	19158	4.571	ng/uL	0.00
4) Pyridine-d5	3.787	84	64843	5.500	ng/ul	0.00
7) Phenol-d5	7.093	99	80081	4.817	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	272831	27.351	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	248246	20.164	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	165573	12.021	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	172817	27.614	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	151906	22.225	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	279951	21.040	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	424440	22.247	ng/ul	-0.01
46) Dimethylphthalate-d6	13.975	166	1231653	29.342	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1359125	26.753	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143	15500m	2.006	ng/ul	0.00
60) Fluorene-d10	15.557	176	1062805	29.434	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	9275m	1.273	ng/ul	0.00
73) Anthracene-d10	17.416	188	1834862	32.904	ng/ul	0.00
81) Pyrene-d10	19.645	212	2046535	37.538	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1277145	34.670	ng/ul	0.00

Target Compounds Qvalue

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

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