

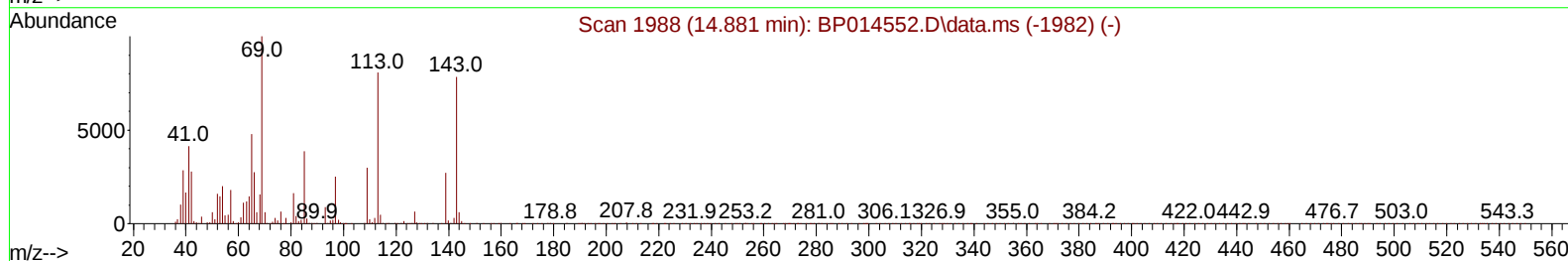
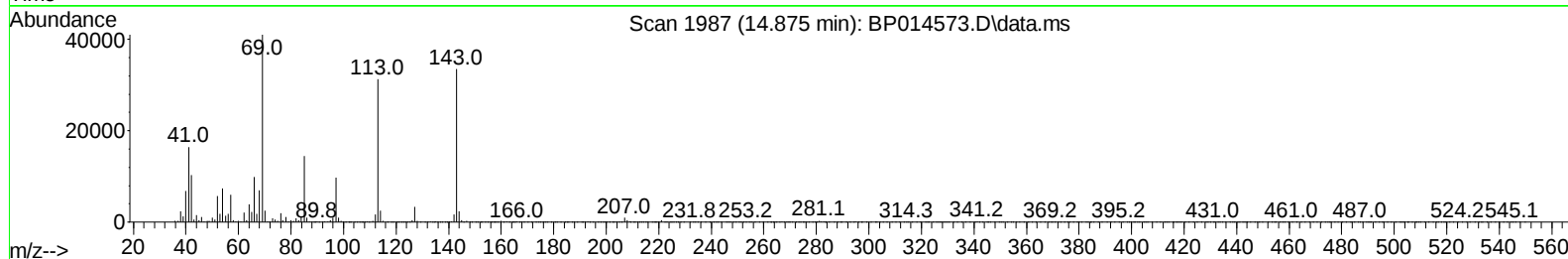
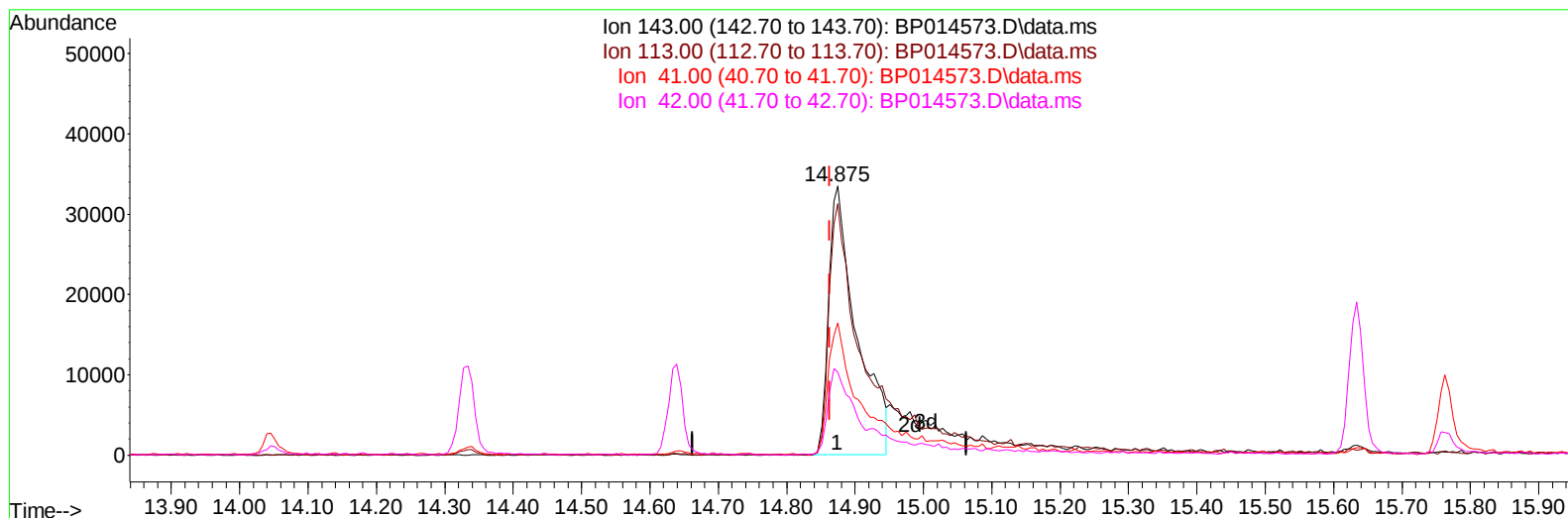
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014573.D
Acq On : 06 Apr 2023 00:28
Operator : CG/JU
Sample : PB151910BL
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK910

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/06/2023
Supervised By :Jagrut Upadhyay 04/06/2023

Quant Time: Apr 06 02:13:05 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 06 00:55:13 2023
Response via : Initial Calibration



TIC: BP014573.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.012) 18.28 ng/u1

response 94939

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	93.55
41.00	44.20	49.15
42.00	31.50	30.77

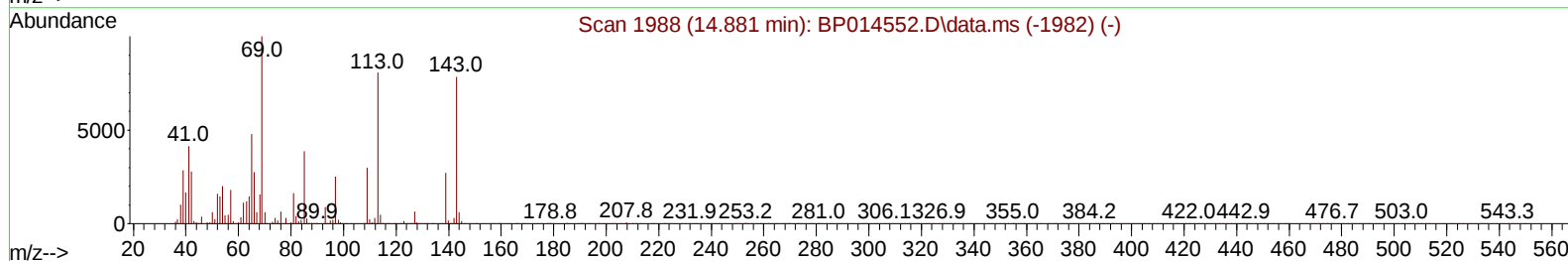
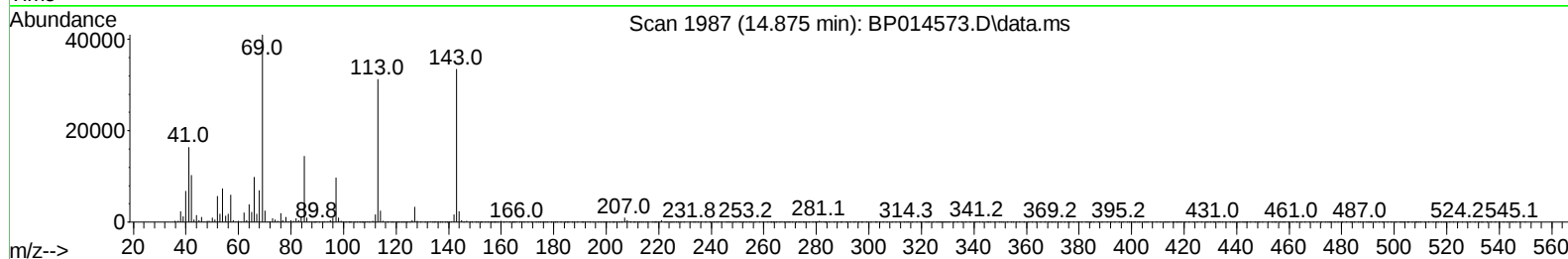
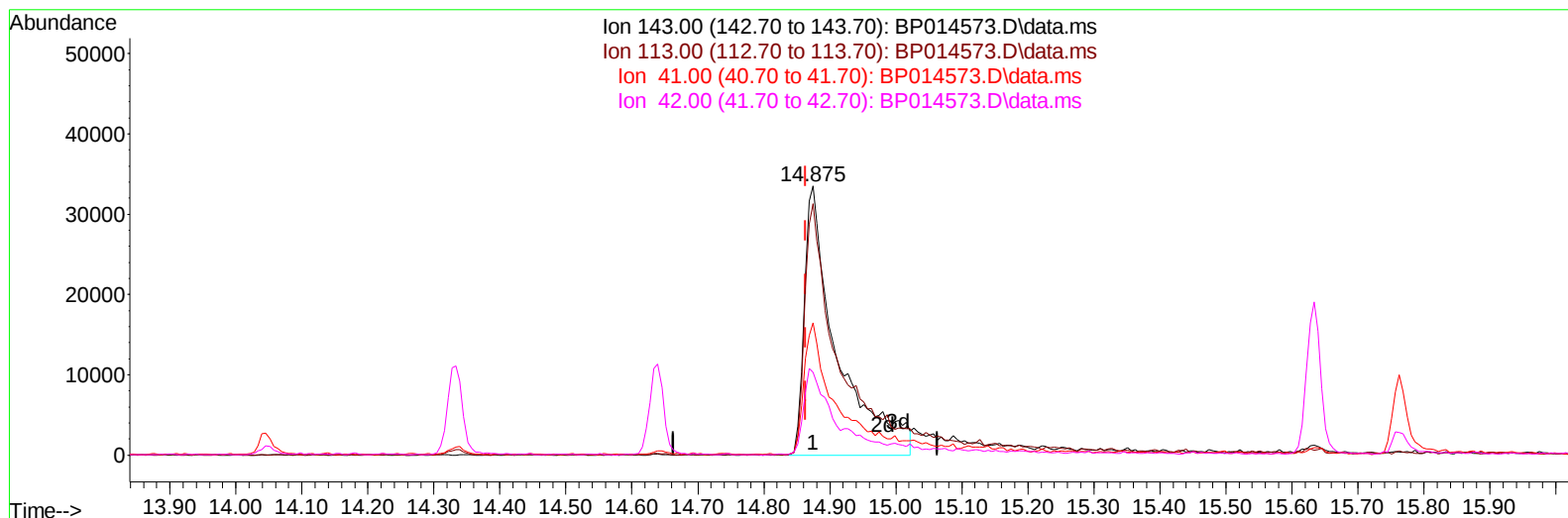
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TIC: BP014573.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.875min (+ 0.012) 22.36 ng/ul m

response 116166

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	93.55
41.00	44.20	49.15
42.00	31.50	30.77

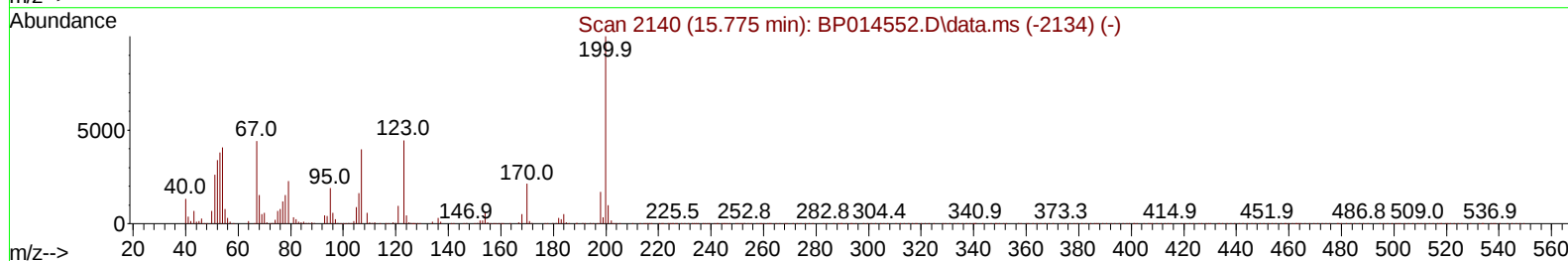
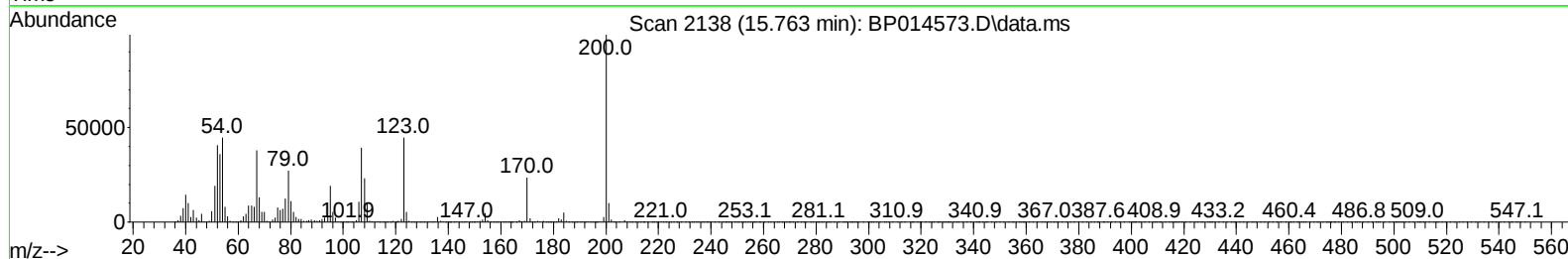
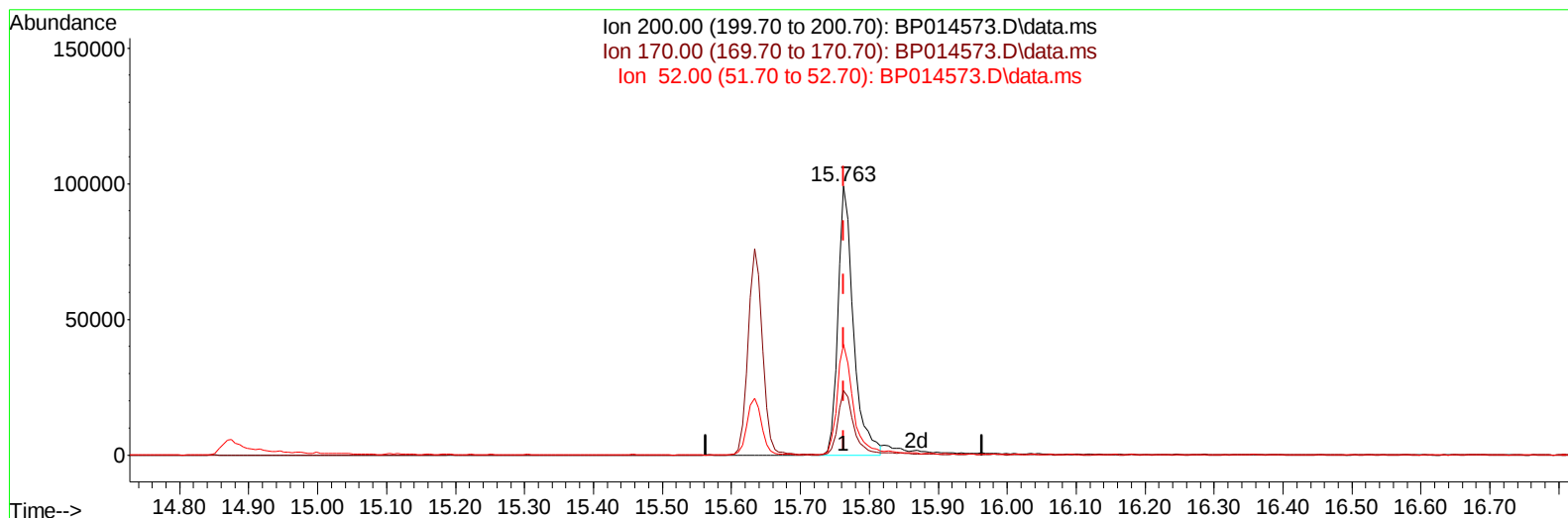
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TIC: BP014573.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.763min (-0.000) 21.43 ng/ul****response 153486**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	23.87
52.00	45.50	41.11
0.00	0.00	0.00

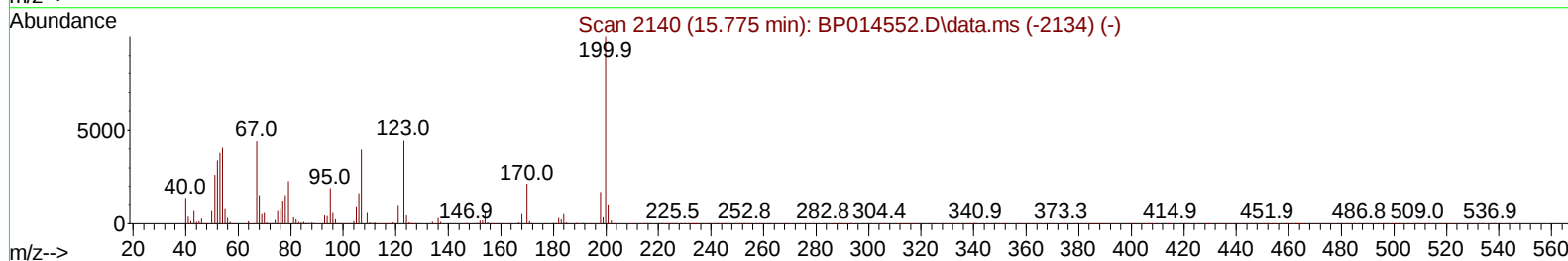
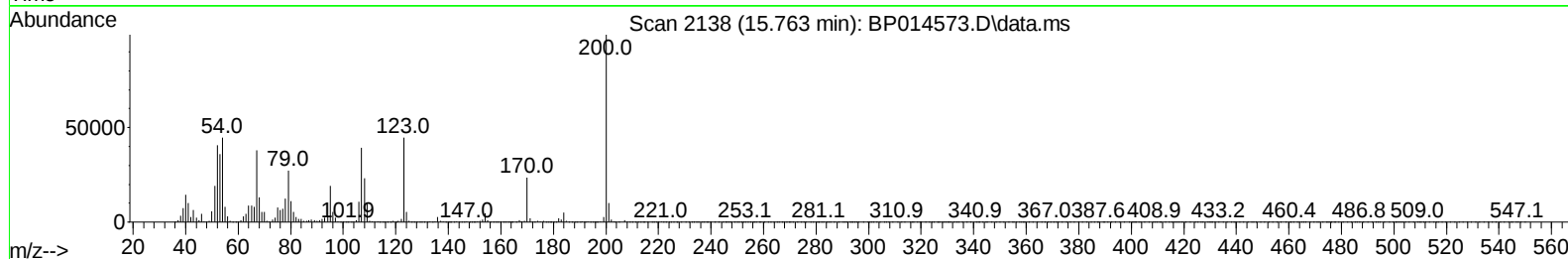
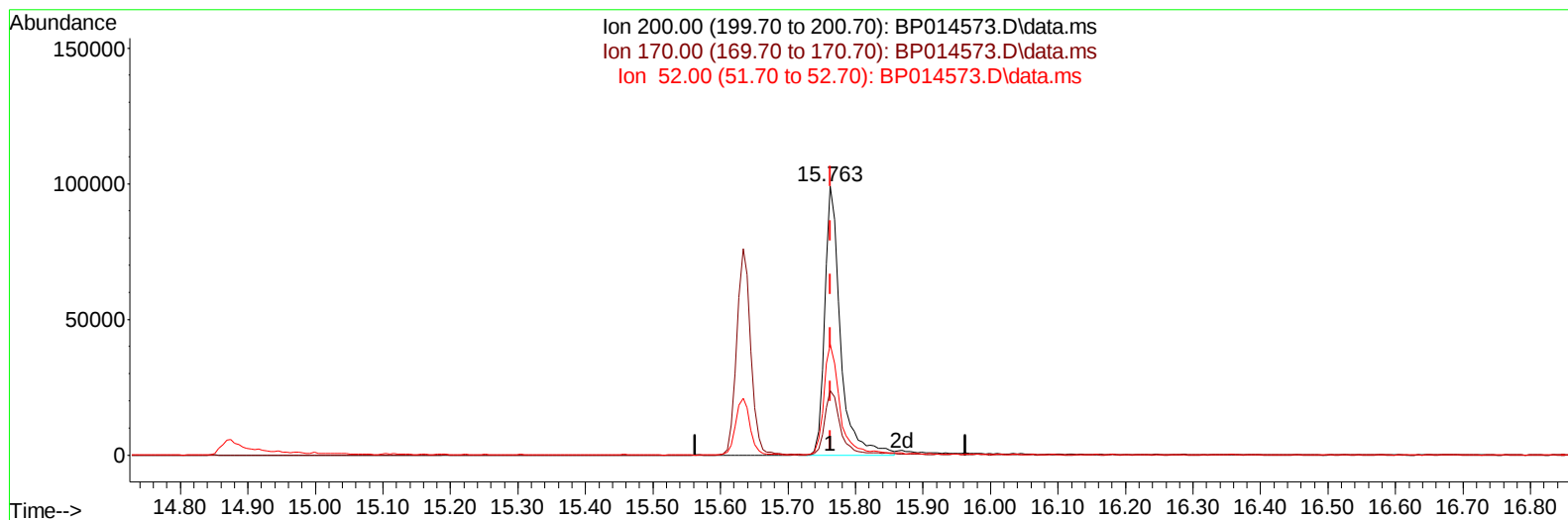
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TIC: BP014573.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.763min (-0.000) 22.31 ng/ul m****response 159786**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.10	23.87
52.00	45.50	41.11
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	7.987	152	140098	20.000	ng/ul	0.00
20) Naphthalene-d8	10.804	136	632365	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.639	164	424082	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	973321	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	886533	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	835394	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	21734	6.016	ng/uL	0.00
4) Pyridine-d5	3.799	84	224883	24.290	ng/ul	0.00
7) Phenol-d5	7.152	99	366378	28.356	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67	266582	32.158	ng/ul	0.00
11) 2-Chlorophenol-d4	7.516	132	284547	30.470	ng/ul	0.00
15) 4-Methylphenol-d8	8.704	113	301995	28.976	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	146721	28.741	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143	161362	27.604	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	270258	25.386	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	497692	35.492	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	1099164	31.966	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	1152371	30.203	ng/ul	0.00
54) 4-Nitrophenol-d4	14.875	143	116166m	22.362	ng/ul	0.01
60) Fluorene-d10	15.633	176	897948	31.315	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.763	200	159786m	22.311	ng/ul	0.00
73) Anthracene-d10	17.498	188	1452610	30.838	ng/ul	0.00
81) Pyrene-d10	19.727	212	1735298	29.141	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1388821	31.291	ng/ul	0.00

Target Compounds Qvalue

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

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