

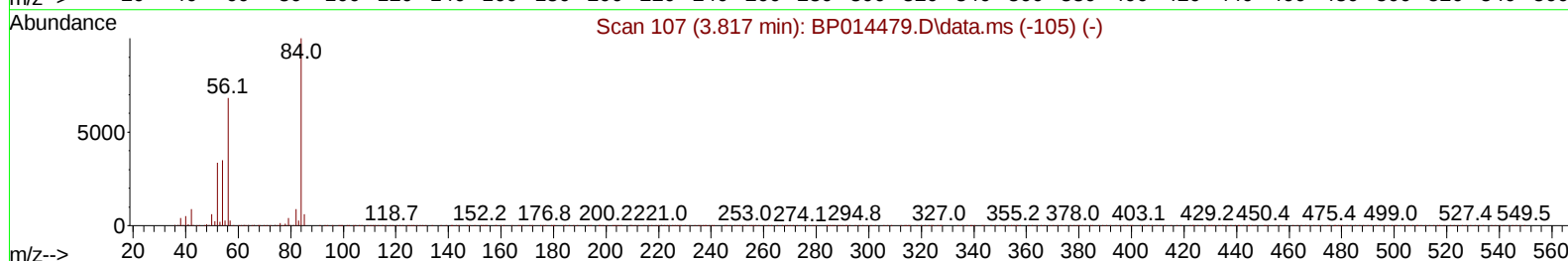
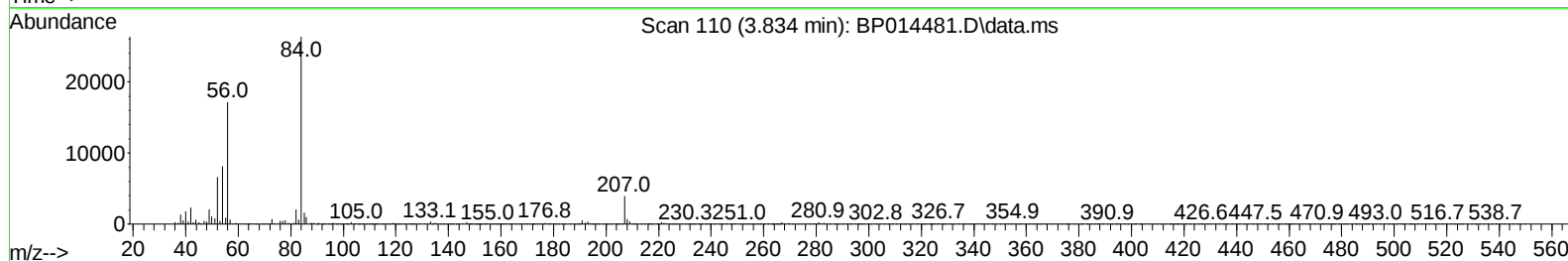
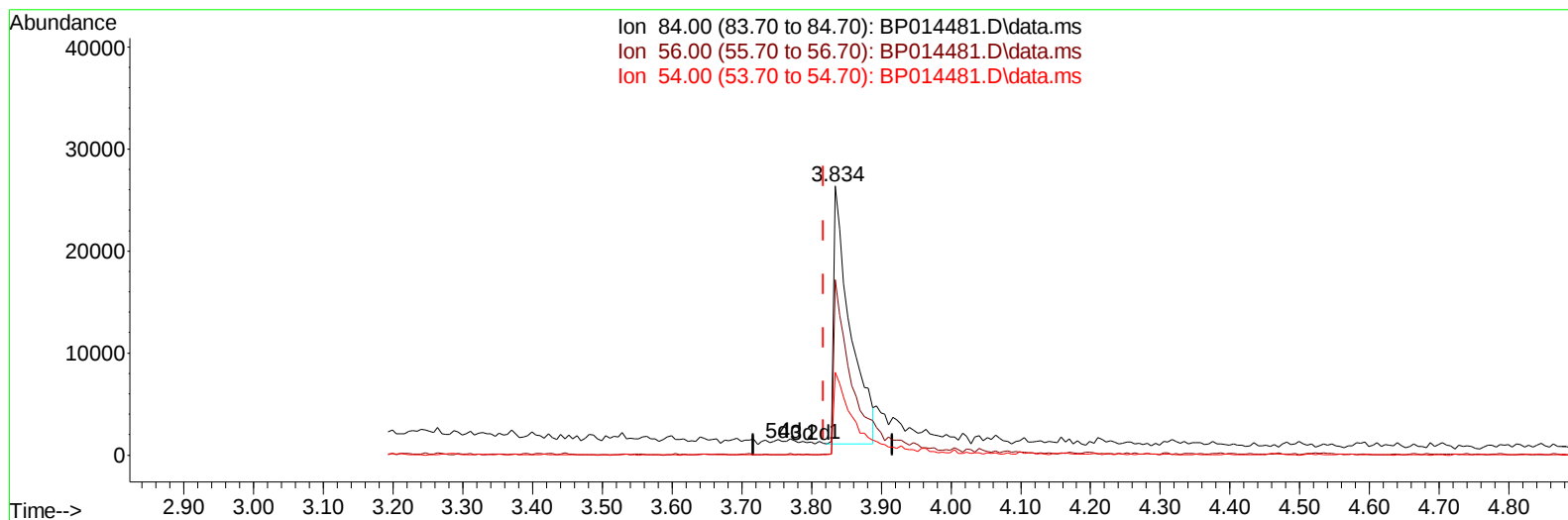
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014481.D
Acq On : 03 Apr 2023 11:51
Operator : CG/JU
Sample : 02158-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-7S(20230329)

Manual IntegrationsAPPROVED

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

Quant Time: Apr 04 00:51:47 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration



TIC: BP014481.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 3.96 ng/u1

response 40566

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.25
54.00	32.60	30.69
0.00	0.00	0.00

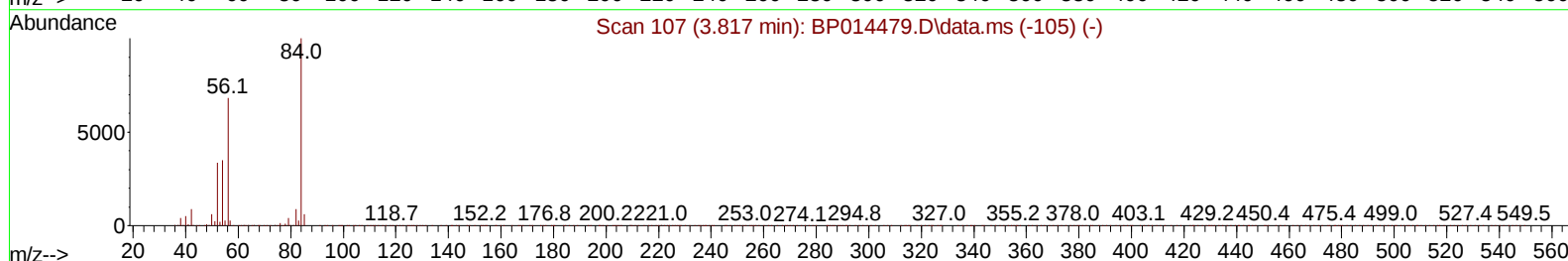
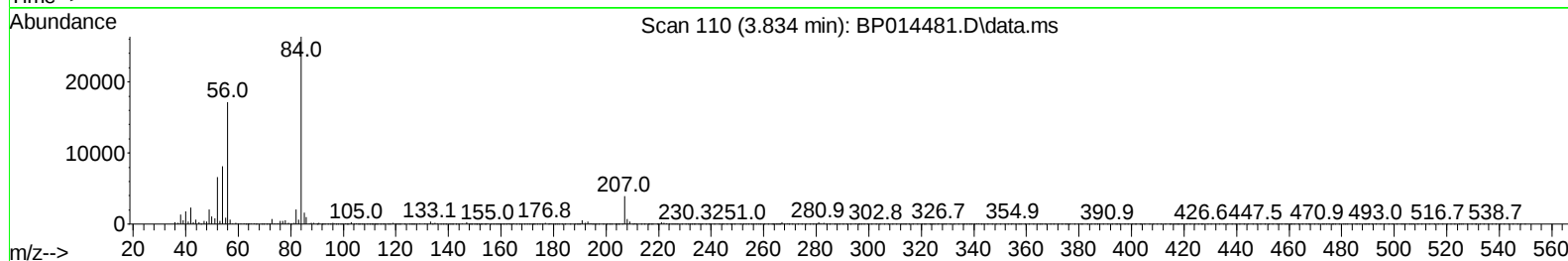
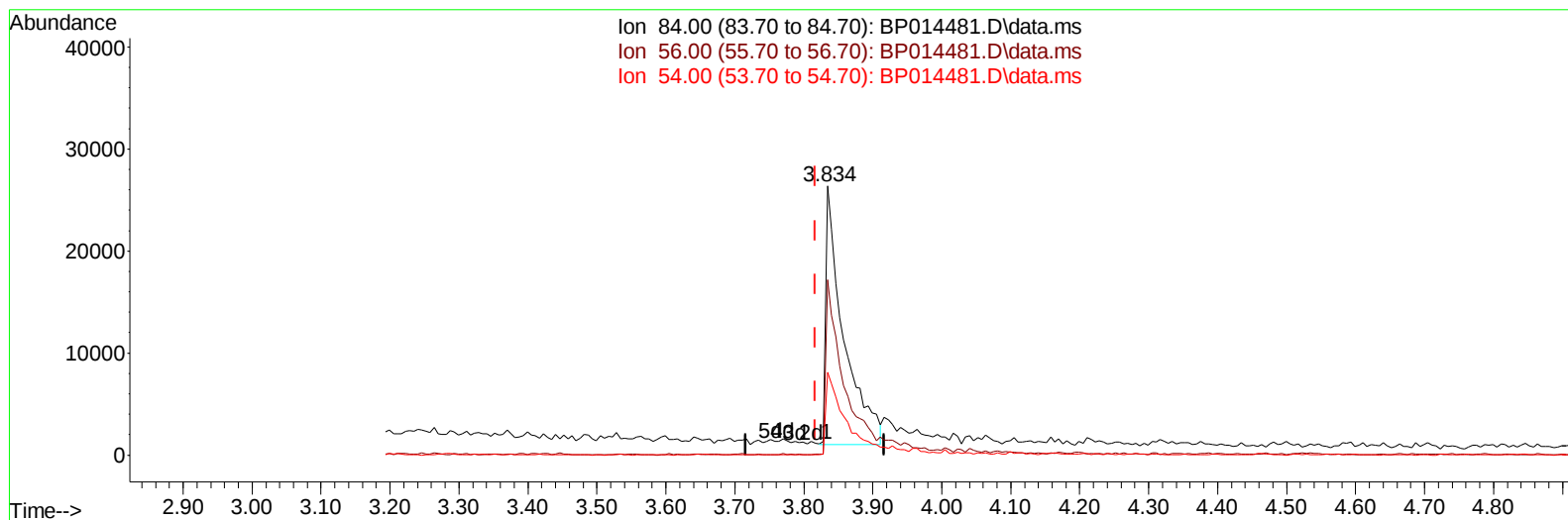
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TIC: BP014481.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 4.38 ng/ul m

response 44901

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.25
54.00	32.60	30.69
0.00	0.00	0.00

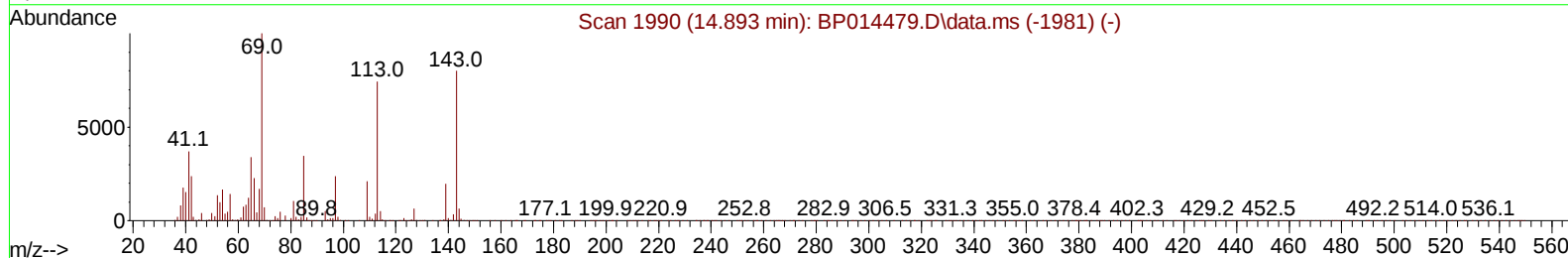
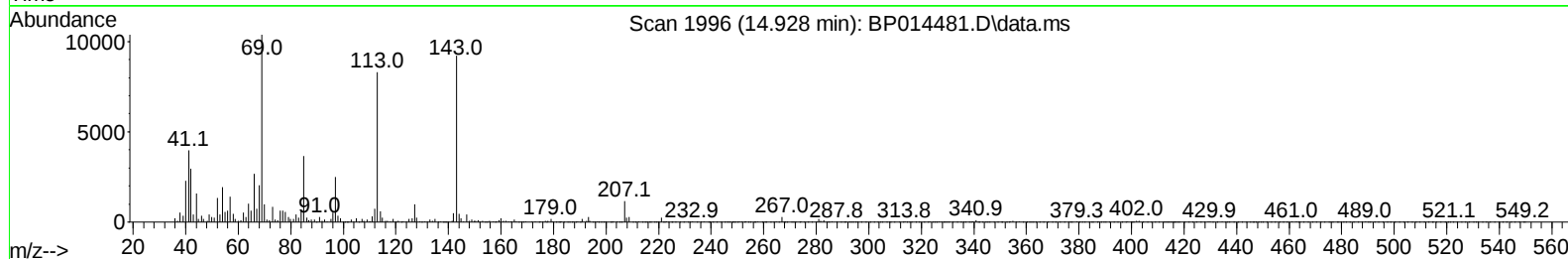
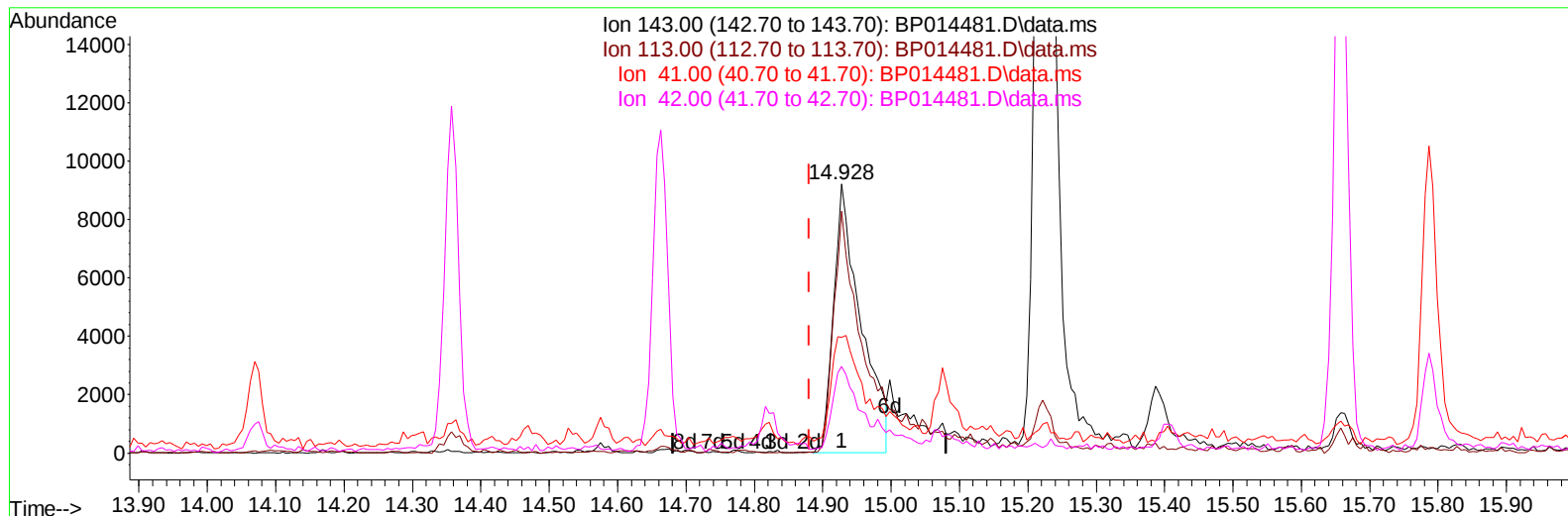
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TIC: BP014481.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.047) 4.45 ng/u1

response 25369

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	90.03
41.00	44.20	43.13
42.00	31.50	32.10

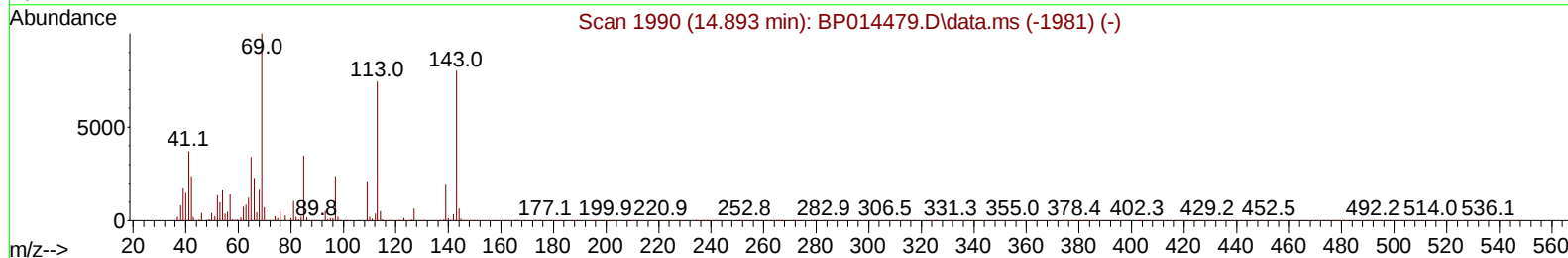
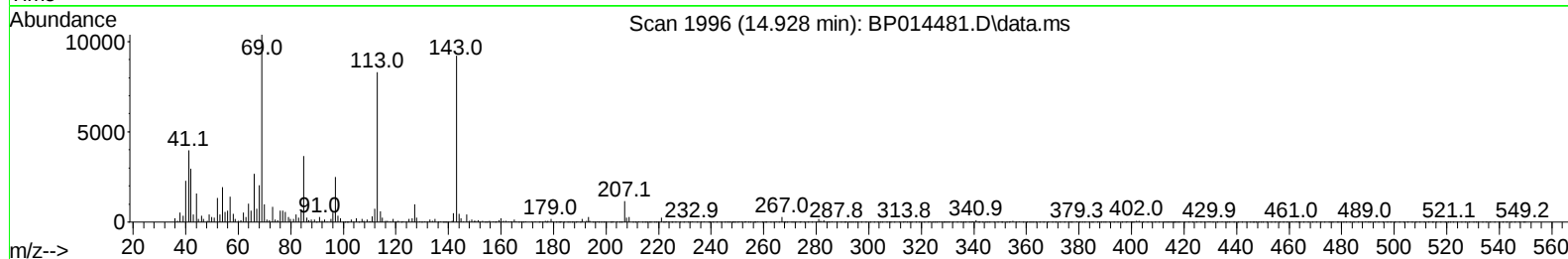
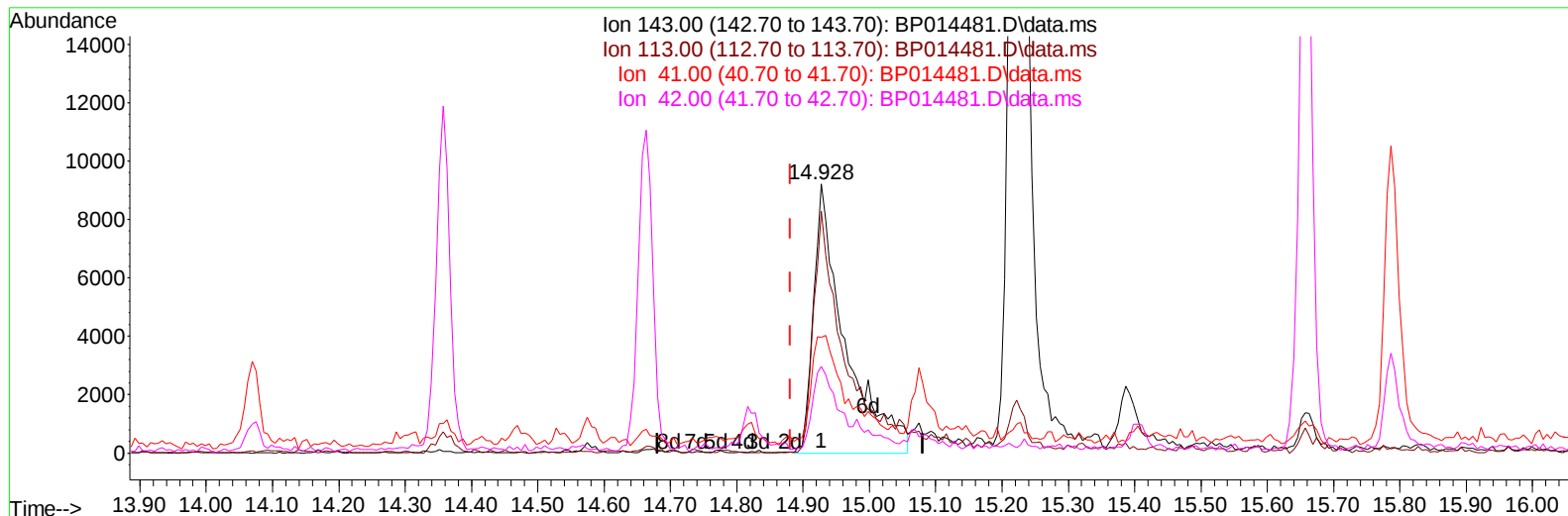
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TIC: BP014481.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.928min (+ 0.047) 5.35 ng/ul m

response 30454

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	90.03
41.00	44.20	43.13
42.00	31.50	32.10

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	154973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	682978	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.663	164	465029	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1062876	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	940726	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	907140	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	16988	4.251	ng/uL	0.00
4) Pyridine-d5	3.834	84	44901m	4.384	ng/ul	0.02
7) Phenol-d5	7.181	99	92445	6.468	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	258934	28.237	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	232632	22.520	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	177295	15.379	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	146965	26.655	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	166534	26.378	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.463	165	287007	24.962	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	261993	17.299	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	1152109	30.556	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1217103	29.091	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	30454m	5.346	ng/ul	0.05
60) Fluorene-d10	15.657	176	971055	30.882	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.787	200	172075	22.002	ng/ul	0.00
73) Anthracene-d10	17.528	188	1596501	31.037	ng/ul	0.00
81) Pyrene-d10	19.757	212	1935624	30.633	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.869	264	1562824	32.427	ng/ul	0.00

Target Compounds Qvalue

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

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