

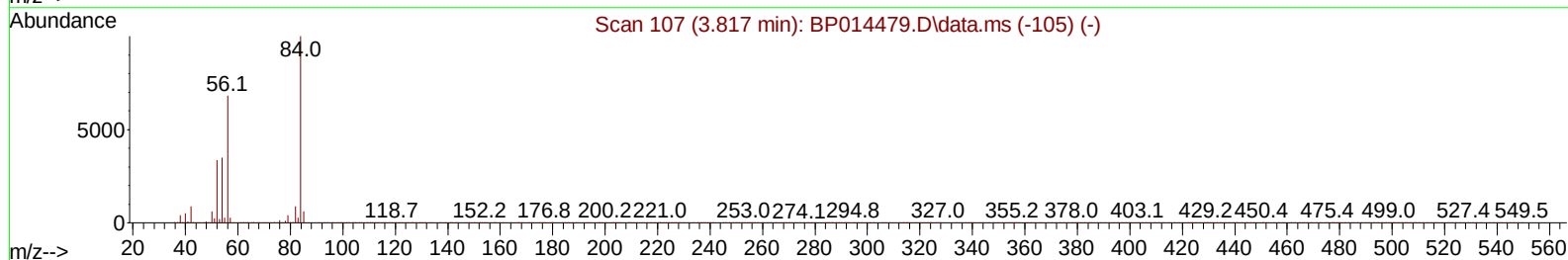
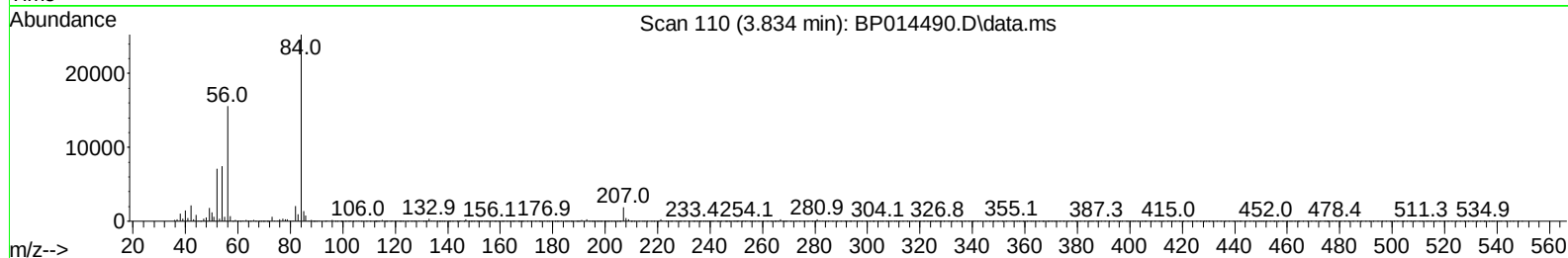
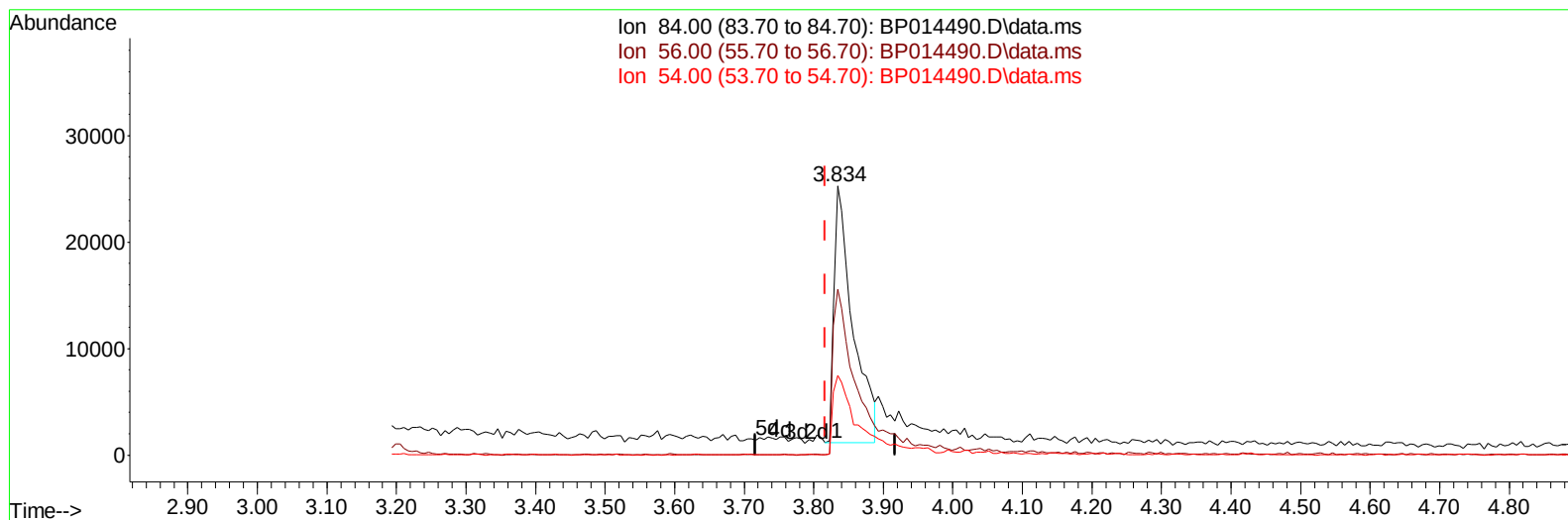
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014490.D
Acq On : 03 Apr 2023 16:57
Operator : CG/JU
Sample : PB151781BL
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK781

Manual IntegrationsAPPROVED

Quant Time: Apr 04 00:54:56 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

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



TIC: BP014490.D\data.ms

(4) Pyridine-d5 (S)

3.834min (+ 0.018) 9.69 ng/uL

response 45016

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	61.61
54.00	32.60	29.64
0.00	0.00	0.00

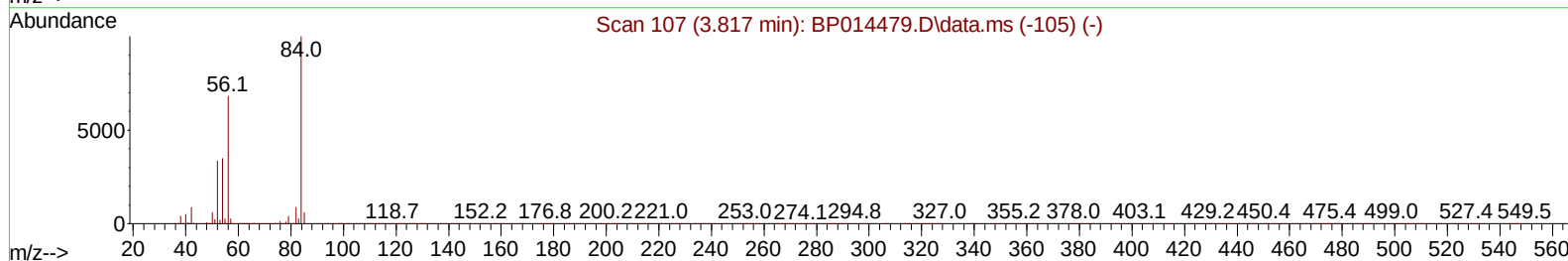
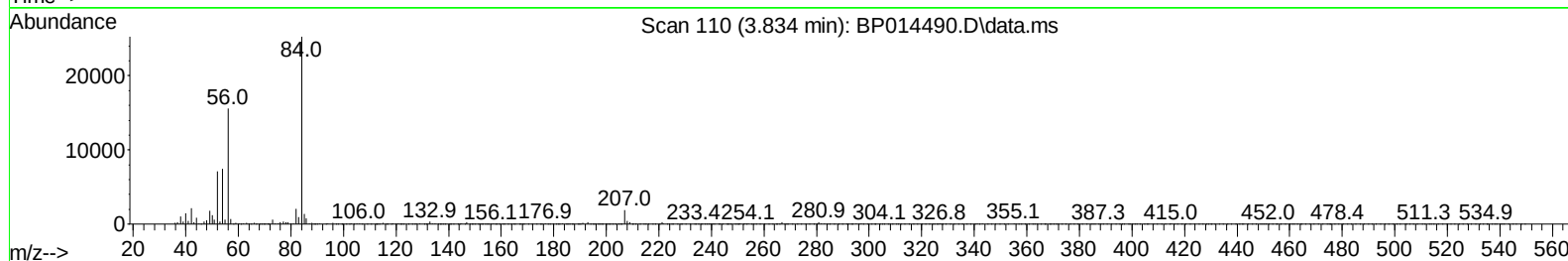
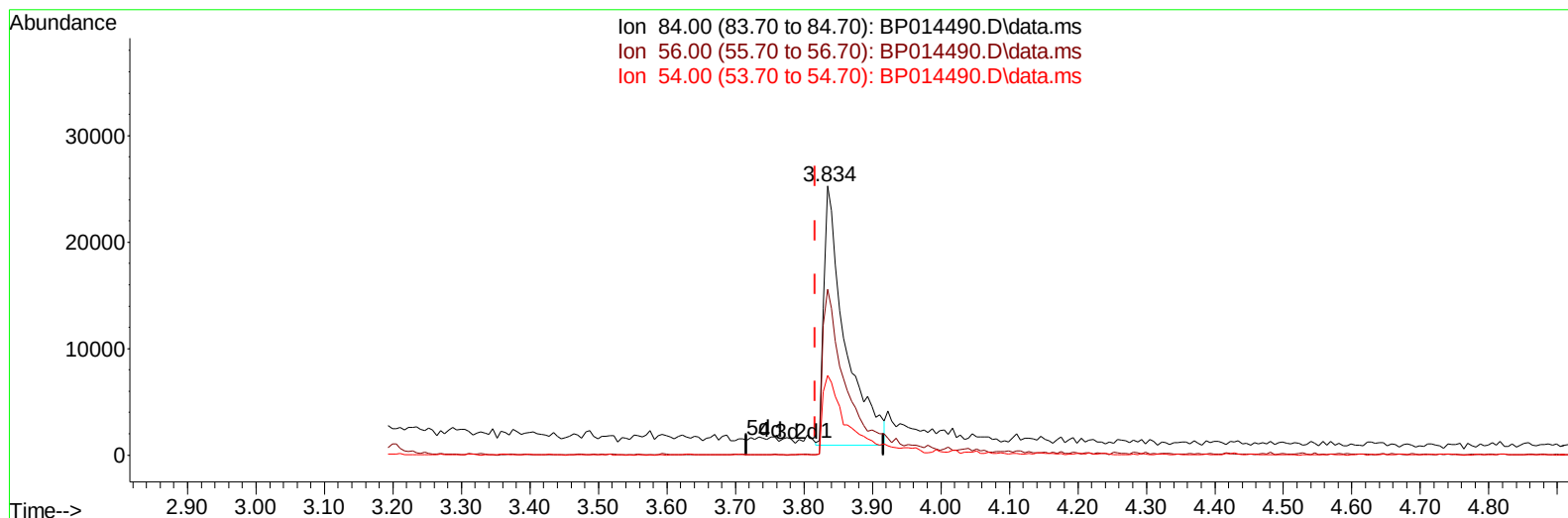
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(4) Pyridine-d5 (S)

3.834min (+ 0.018) 11.14 ng/ul m

response 51745

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	61.61
54.00	32.60	29.64
0.00	0.00	0.00

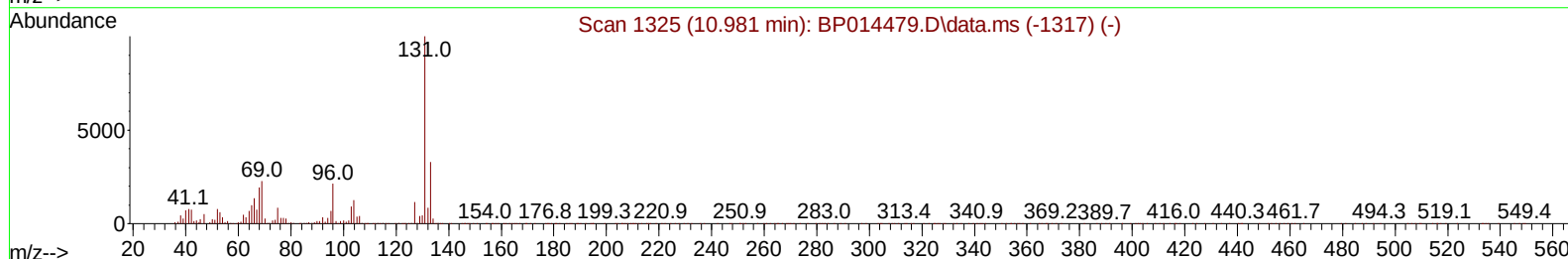
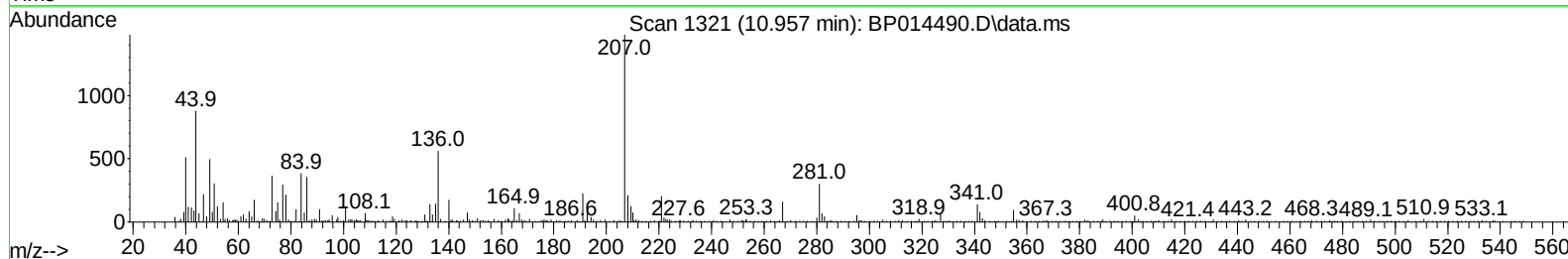
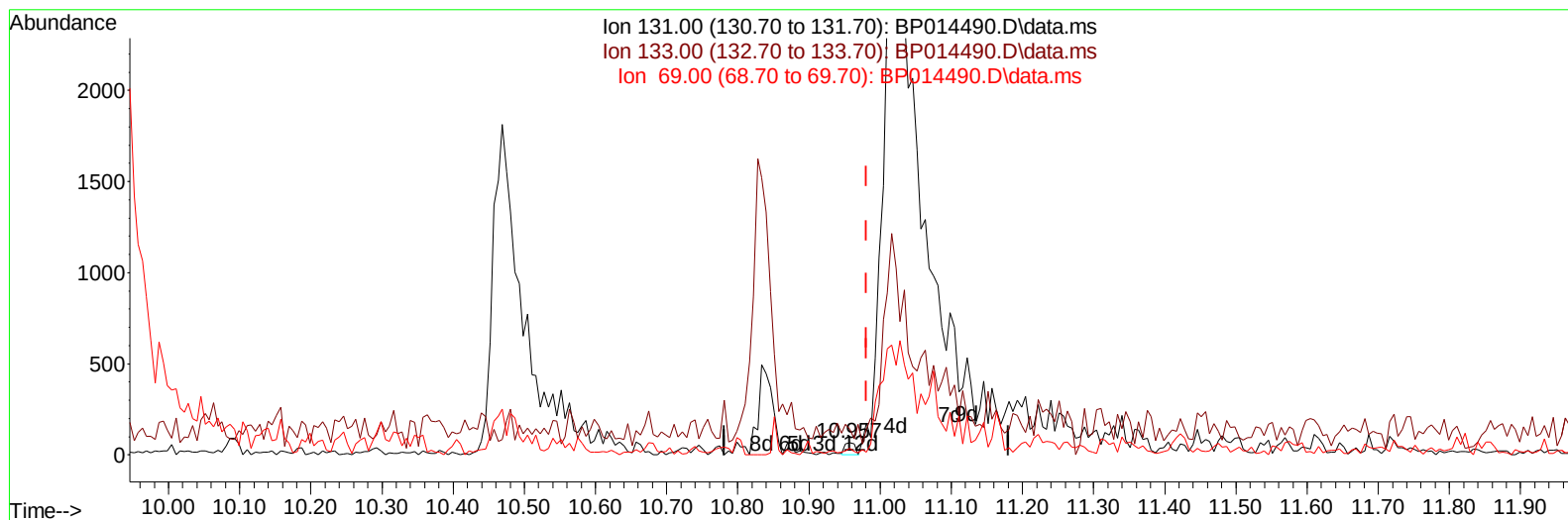
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TIC: BP014490.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.957min (-0.024) 0.01 ng/u1

response 67

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	211.94#
69.00	21.30	55.22#
0.00	0.00	0.00

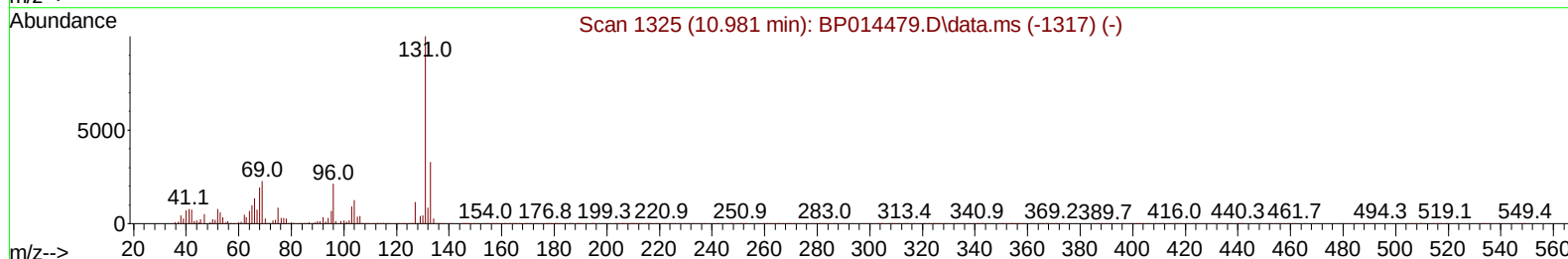
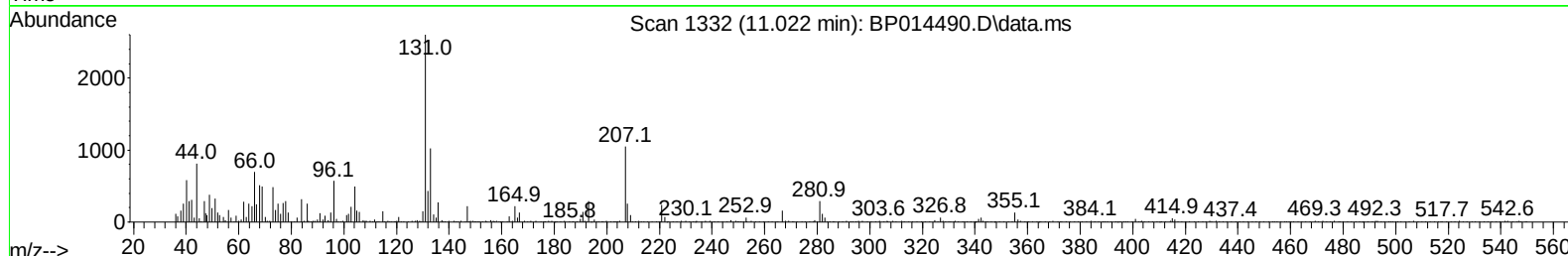
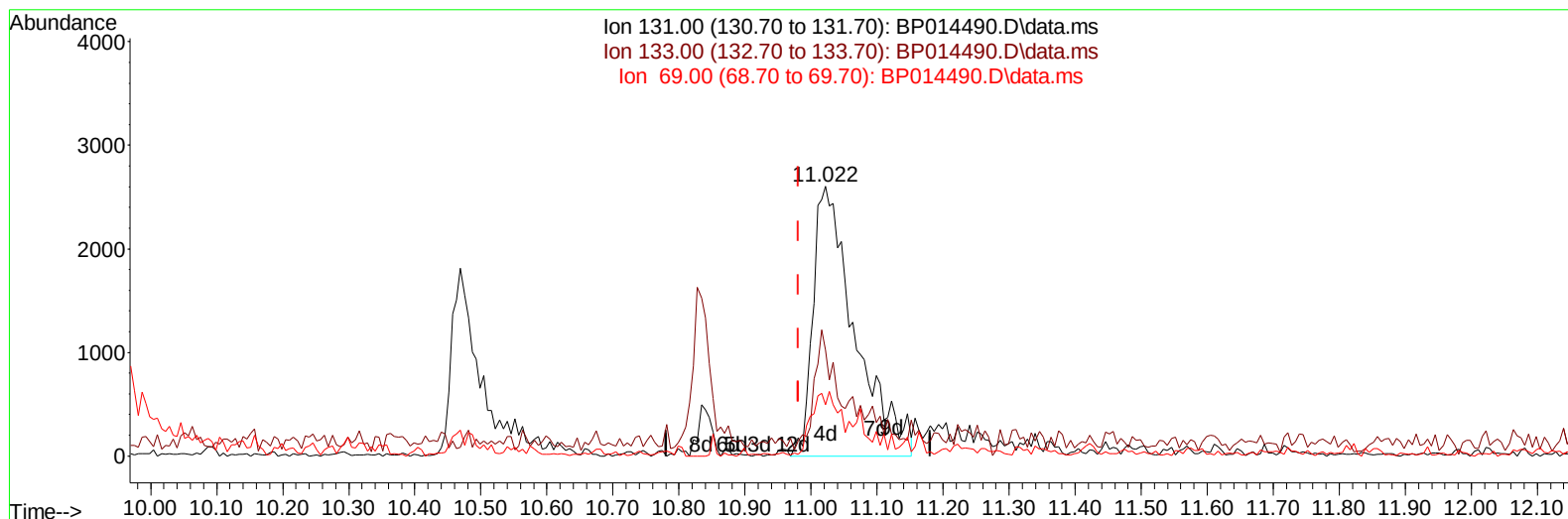
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TIC: BP014490.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.022min (+ 0.041) 1.61 ng/ul m

response 11431

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	39.32#
69.00	21.30	18.91
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	8.005	152	70292	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	319916	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.663	164	228859	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.422	188	590097	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	747343	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	813936	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	9198	5.074	ng/uL	0.00
4) Pyridine-d5	3.834	84	51745m	11.140	ng/ul	0.02
7) Phenol-d5	7.181	99	182582	28.165	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	116963	28.121	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	133407	28.473	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	148090	28.320	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	64983	25.162	ng/ul	0.00
24) 2-Nitrophenol-d4	9.910	143	79183	26.775	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	123296	22.893	ng/ul	0.01
31) 4-Chloroaniline-d4	11.022	131	11431m	1.611	ng/ul	0.04
46) Dimethylphthalate-d6	14.069	166	523602	28.217	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	566735	27.525	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	76377	27.245	ng/ul	0.03
60) Fluorene-d10	15.657	176	471876	30.494	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	85459	19.682	ng/ul	0.00
73) Anthracene-d10	17.522	188	832205	29.141	ng/ul	0.00
81) Pyrene-d10	19.751	212	1151076	22.930	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.863	264	1272318	29.422	ng/ul	0.00

Target Compounds Qvalue

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

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