

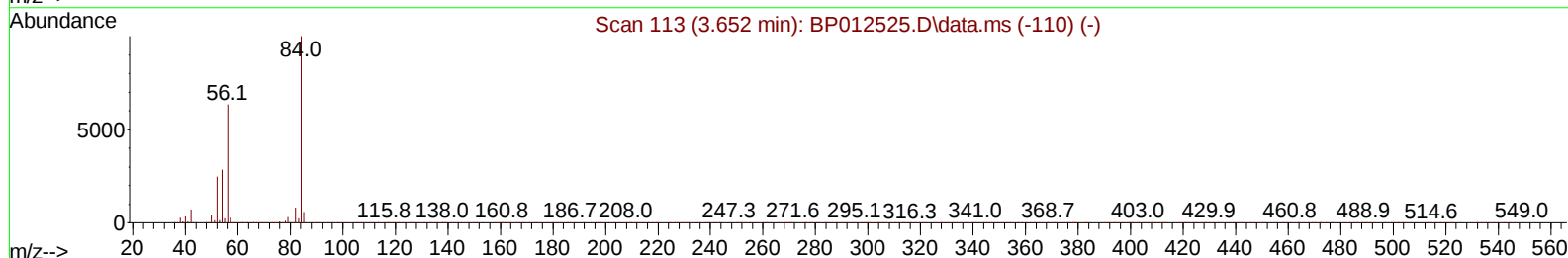
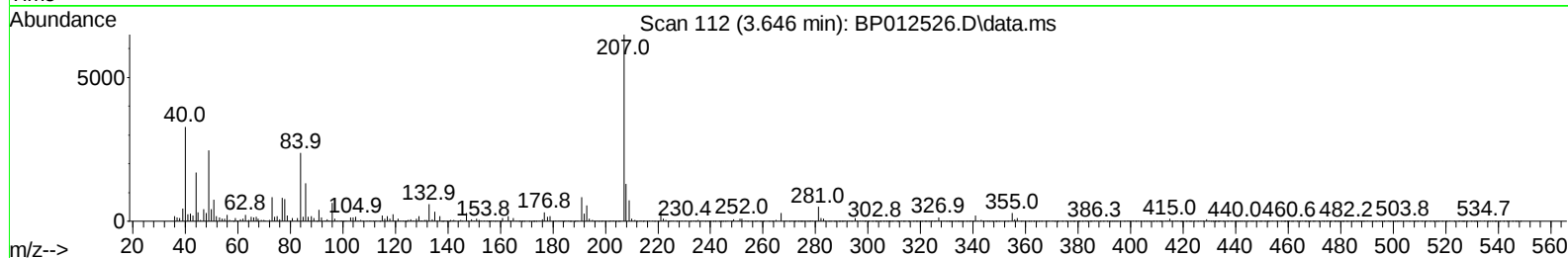
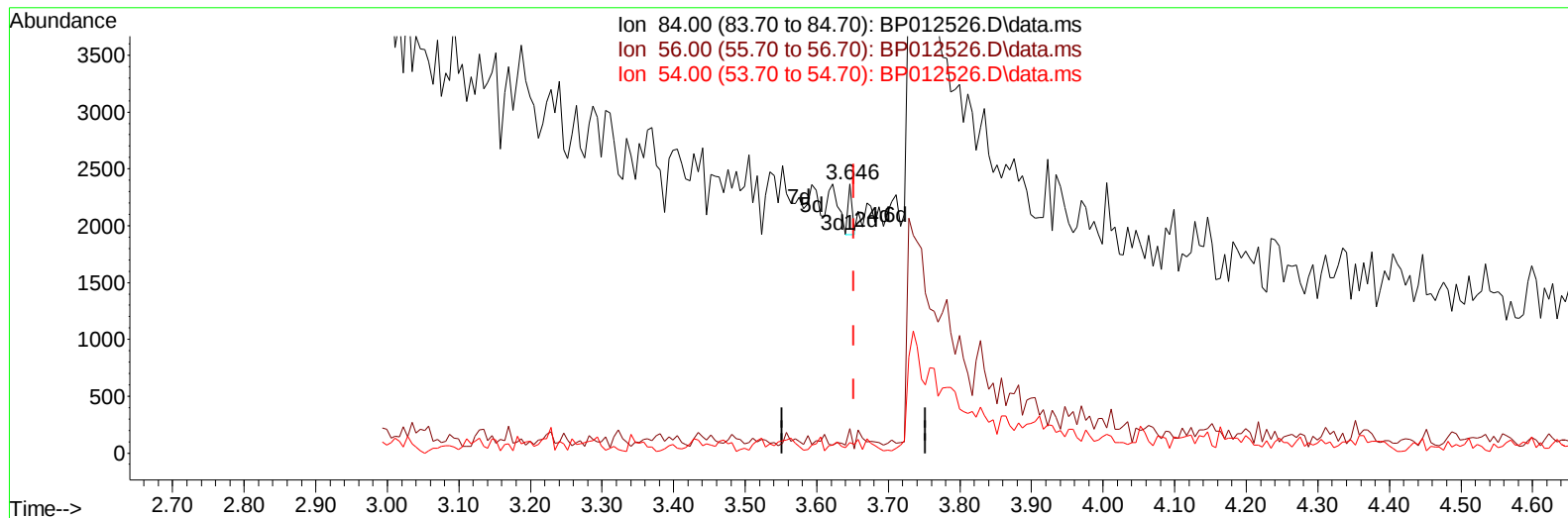
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110722\
Data File : BP012526.D
Acq On : 07 Nov 2022 12:04
Operator : CG/JU
Sample : PB148749BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK749

Manual IntegrationsAPPROVED

Quant Time: Nov 07 21:39:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 04 02:53:36 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/08/2022
Supervised By :mohammad ahmed 11/14/2022



TIC: BP012526.D\data.ms

(4) Pyridine-d5 (S)

3.646min (-0.006) 0.01 ng/u1

response 167

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	9.09#
54.00	29.80	3.97#
0.00	0.00	0.00

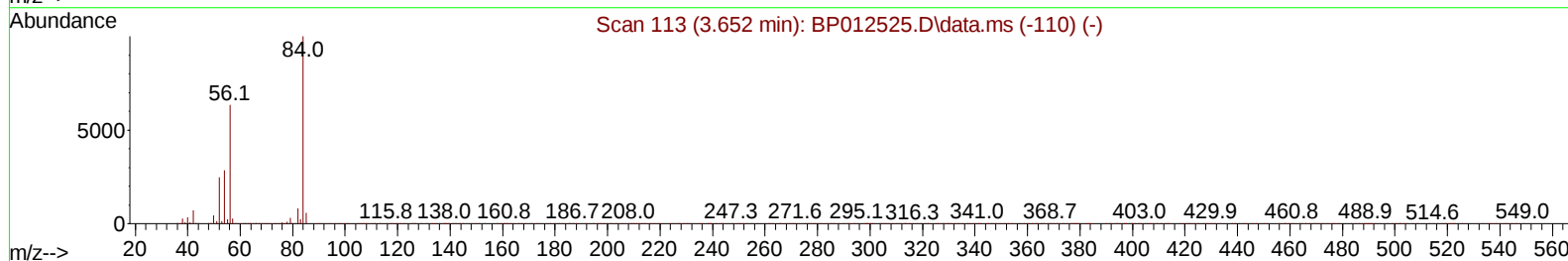
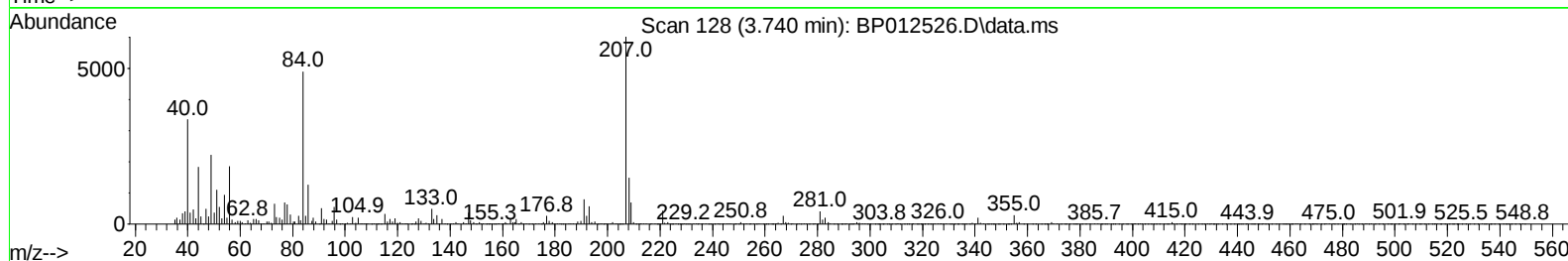
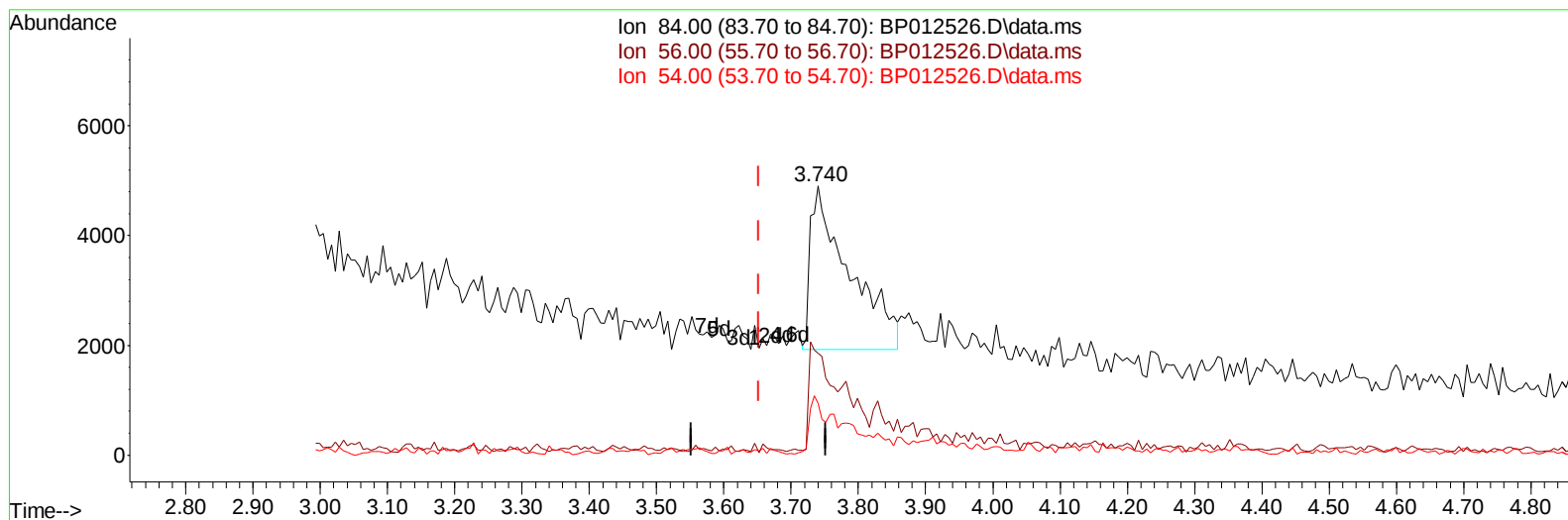
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(4) Pyridine-d5 (S)

3.740min (+ 0.088) 0.61 ng/ul m

response 12054

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	37.94#
54.00	29.80	19.18#
0.00	0.00	0.00

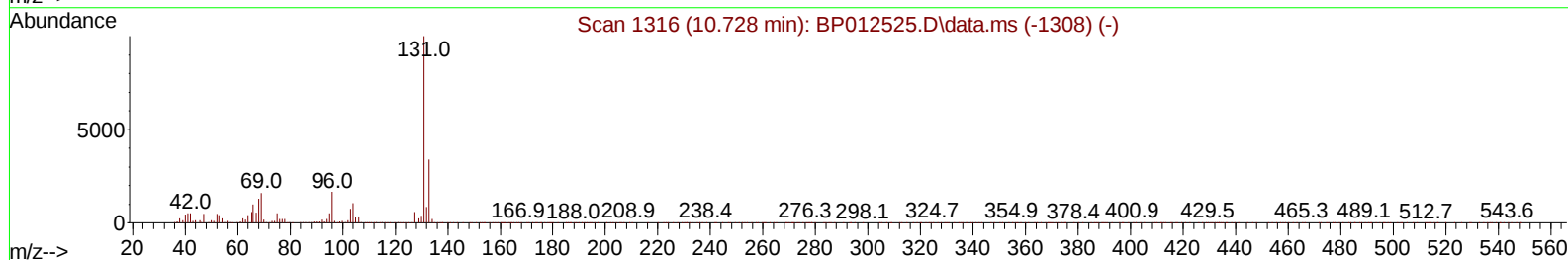
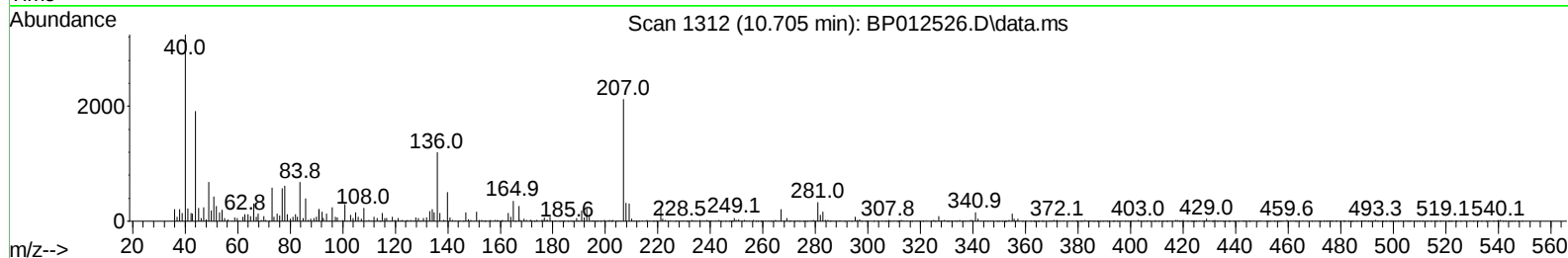
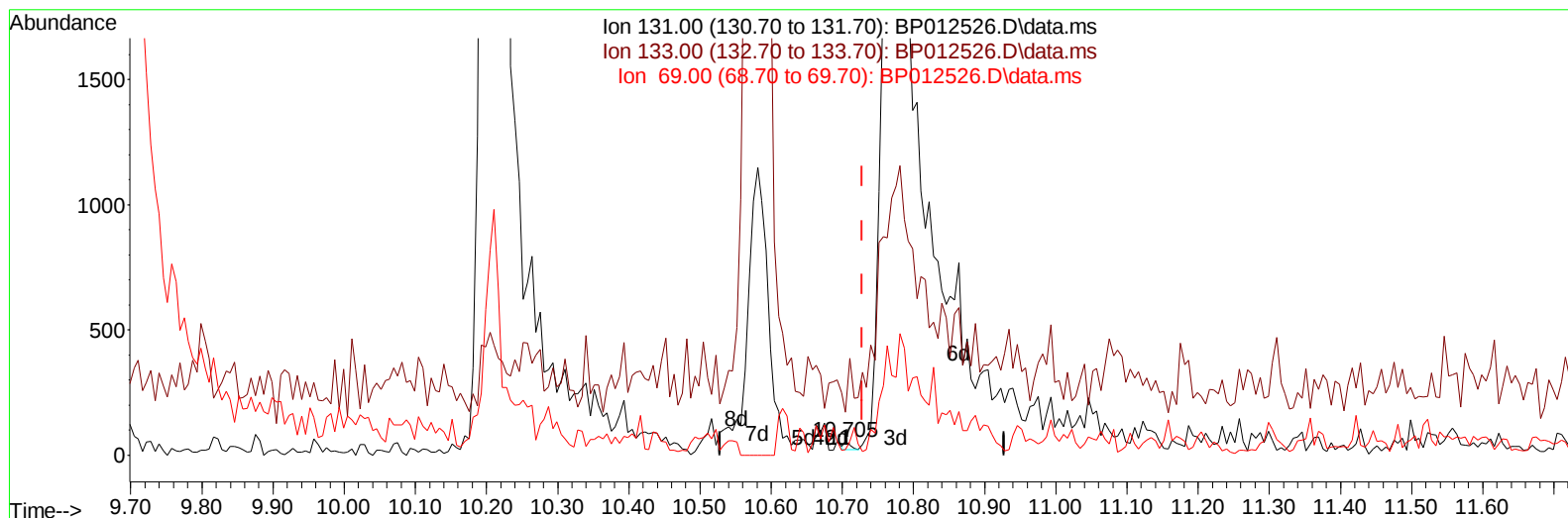
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TIC: BP012526.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.705min (-0.023) 0.00 ng/u1

response 18

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	316.67#
69.00	16.50	40.74#
0.00	0.00	0.00

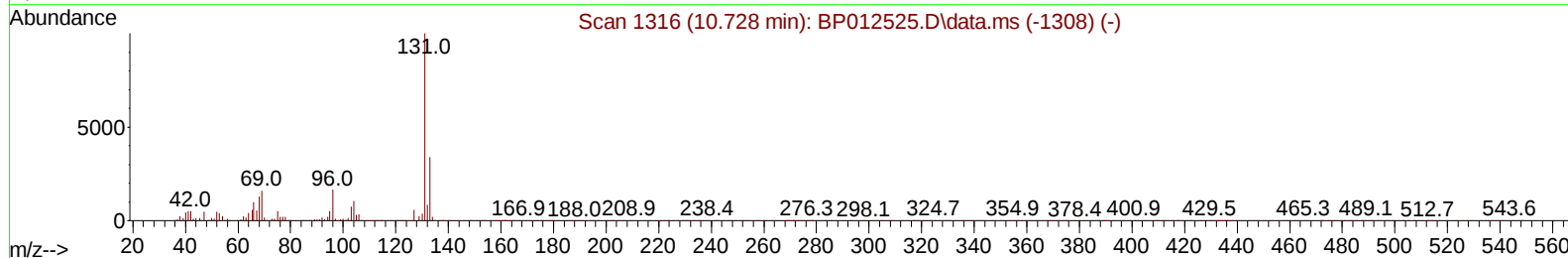
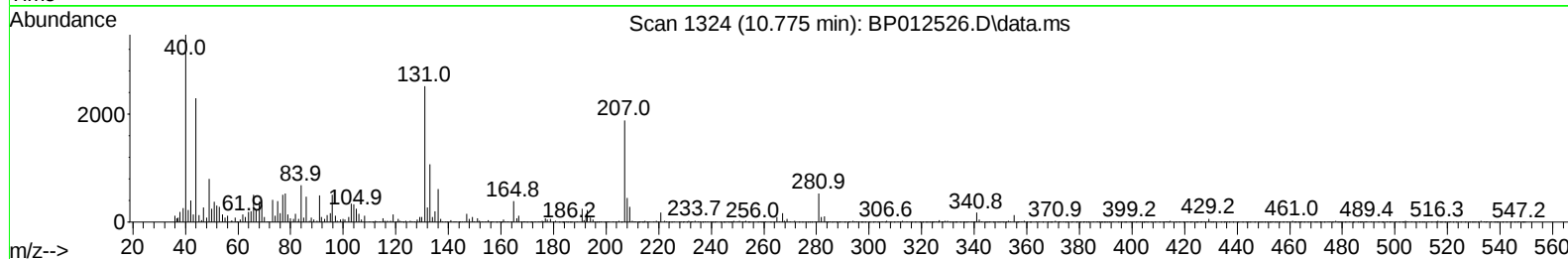
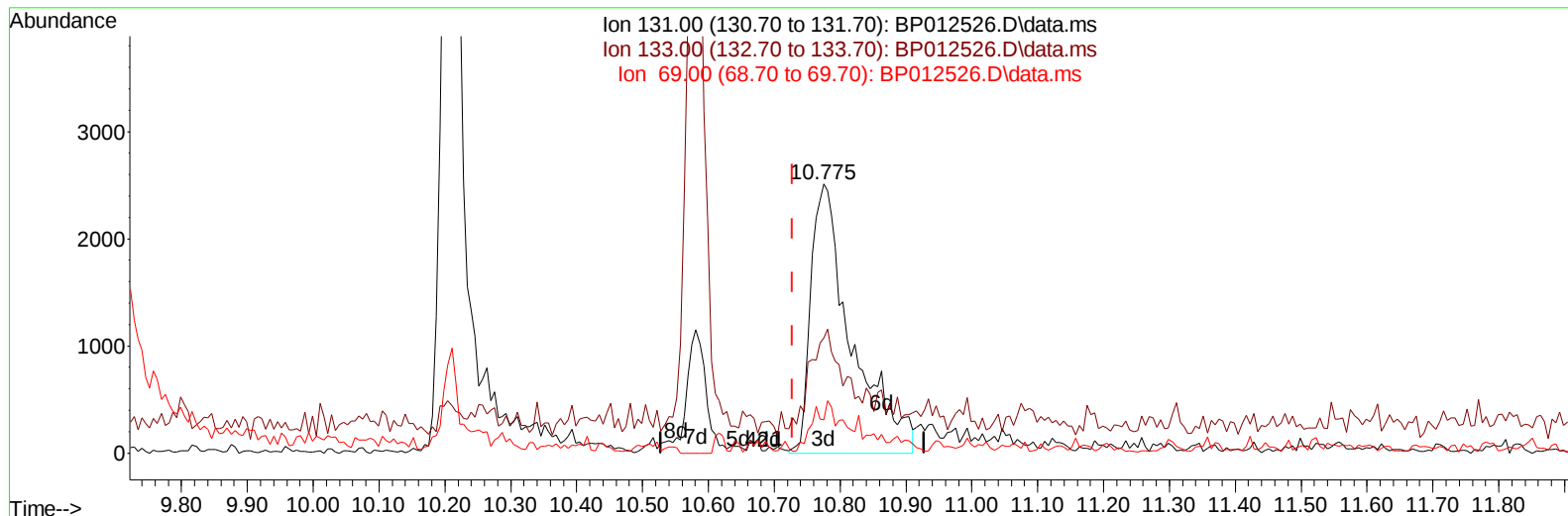
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TIC: BP012526.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.775min (+ 0.047) 0.37 ng/ul m

response 10745

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	42.79#
69.00	16.50	12.43#
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	7.775	152	276803	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	1325077	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	927110	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	2096666	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240	1686717	20.000	ng/ul	0.00
88) Perylene-d12	23.669	264	1482071	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	32828	4.775	ng/uL	0.00
4) Pyridine-d5	3.740	84	12054m	0.610	ng/ul	0.09
7) Phenol-d5	6.958	99	658602	26.528	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	373085	25.173	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	497969	27.466	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	543082	27.658	ng/ul	0.00
21) Nitrobenzene-d5	8.952	128	253377	29.487	ng/ul	0.00
24) 2-Nitrophenol-d4	9.669	143	293432	32.900	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	529201	27.379	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	10745m	0.373	ng/ul	0.05
46) Dimethylphthalate-d6	13.846	166	1775823	27.981	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160	1878287	25.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	213780	17.798	ng/ul	0.01
60) Fluorene-d10	15.428	176	1560324	28.715	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	125575	11.504	ng/ul	0.00
73) Anthracene-d10	17.292	188	2469773	27.338	ng/ul	0.00
81) Pyrene-d10	19.534	212	2716532	32.078	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	2013631	27.806	ng/ul	0.00

Target Compounds Qvalue

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

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