Quantitation Report (Qedit)

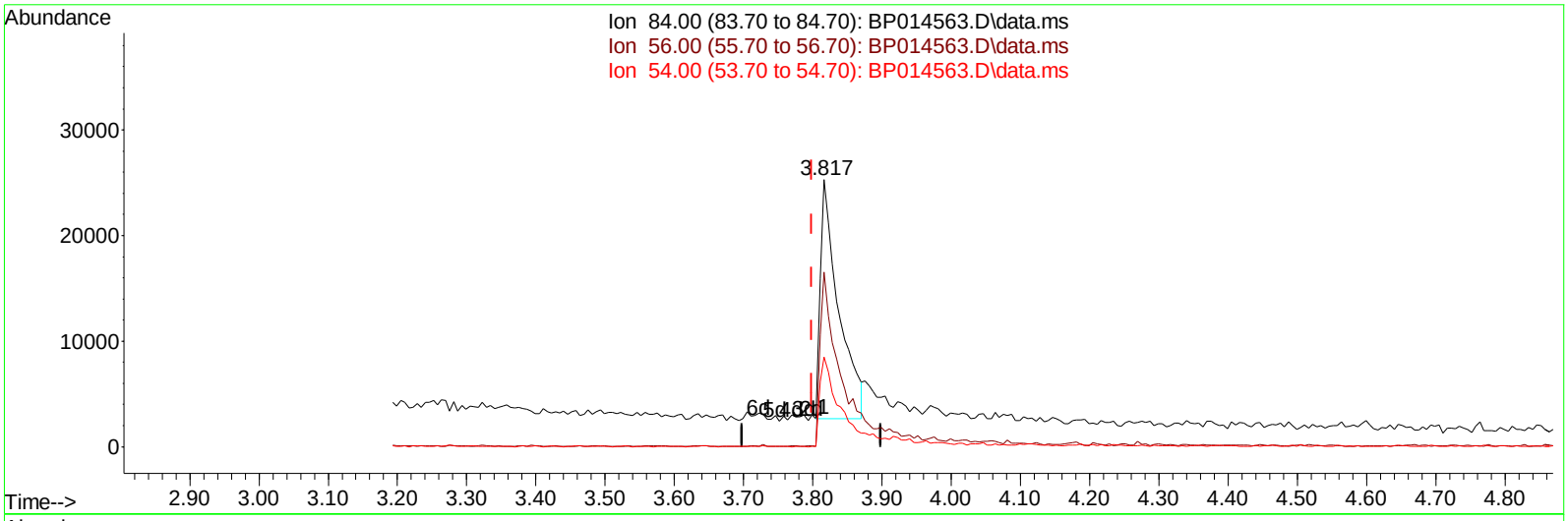
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
Data File : BP014563.D
Acq On : 05 Apr 2023 18:18
Operator : CG/JU
Sample : 02122-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

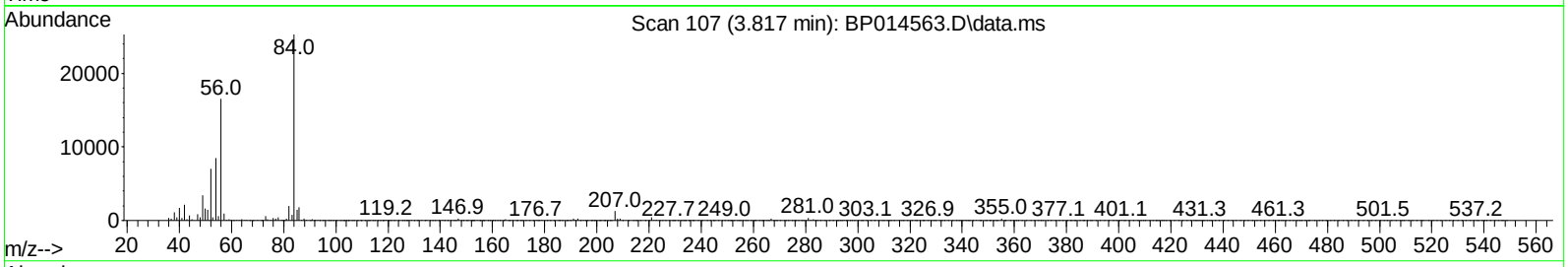
Quant Time: Apr 06 00:58:35 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 06 00:55:13 2023
Response via : Initial Calibration

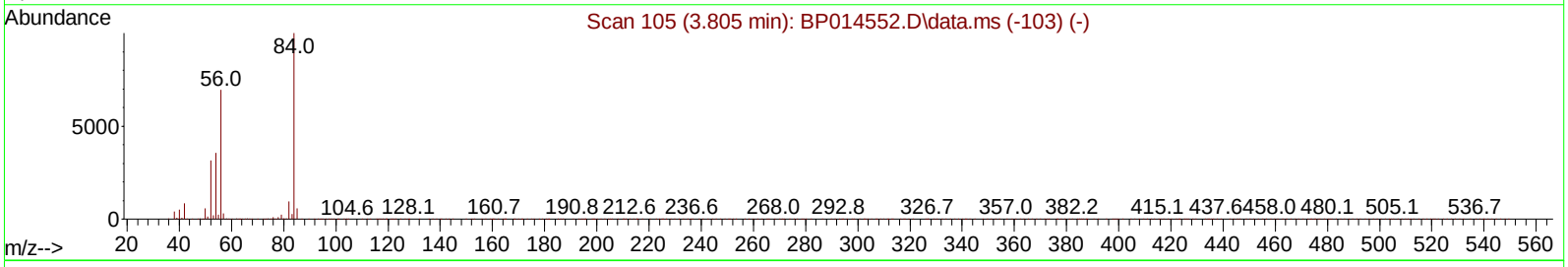
Instrument :
BNA_P
ClientSampleId :
CB9B2

Manual IntegrationsAPPROVED

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







TIC: BP014563.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.018) 5.76 ng/ul

response 41023

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.35
54.00	32.60	33.68
0.00	0.00	0.00

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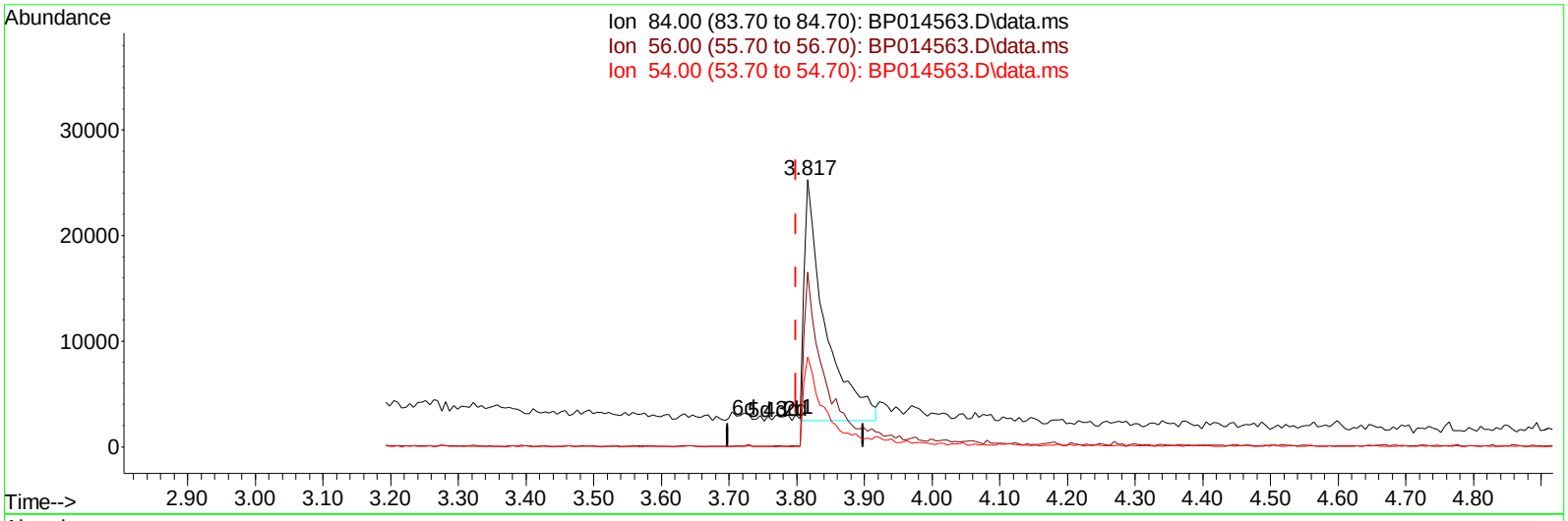

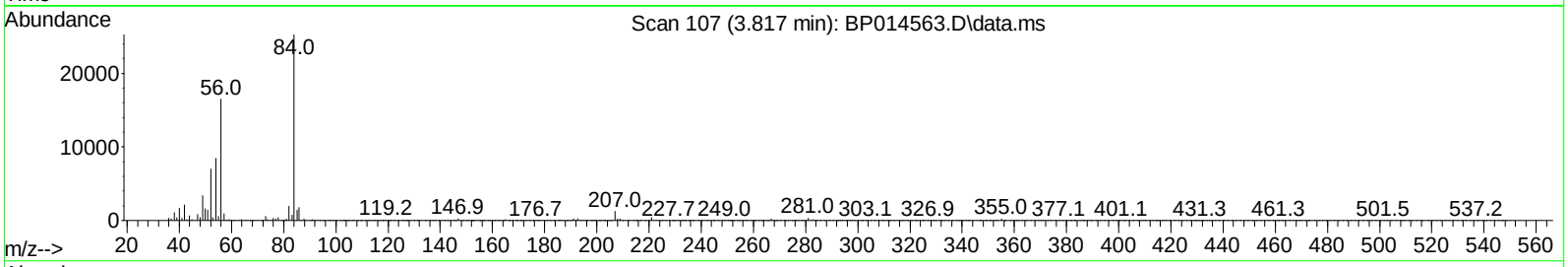

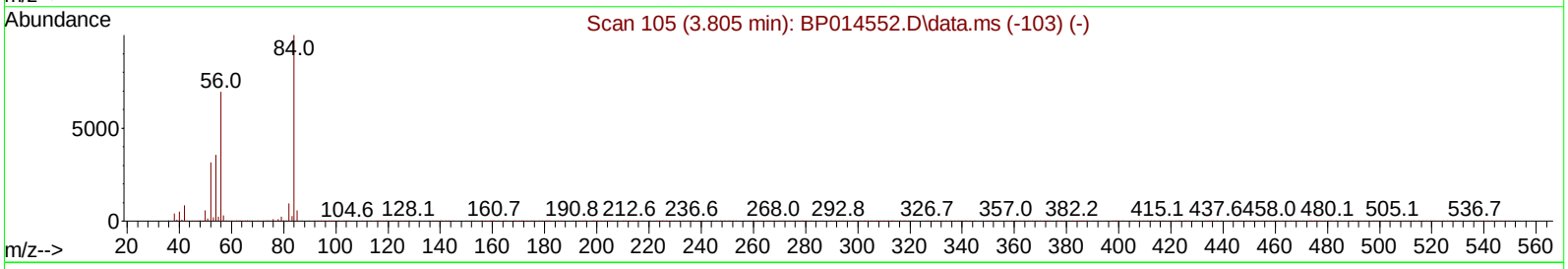

TIC: BP014563.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.018) 6.83 ng/ul m

response 48676

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	65.35
54.00	32.60	33.68
0.00	0.00	0.00

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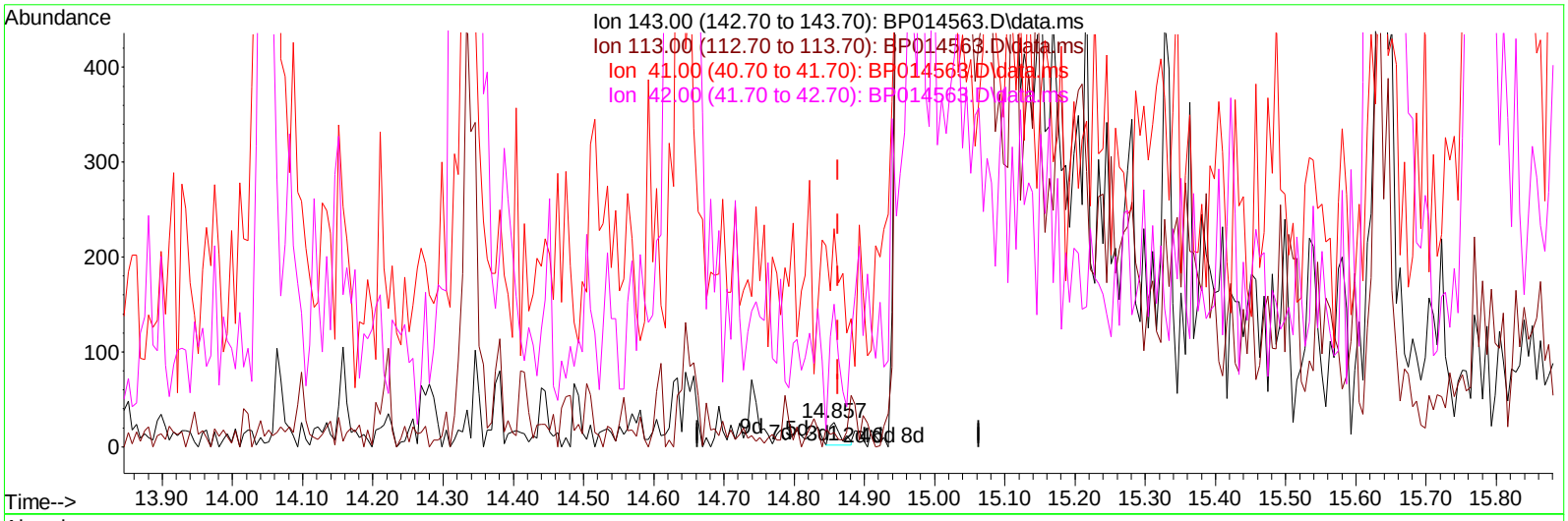

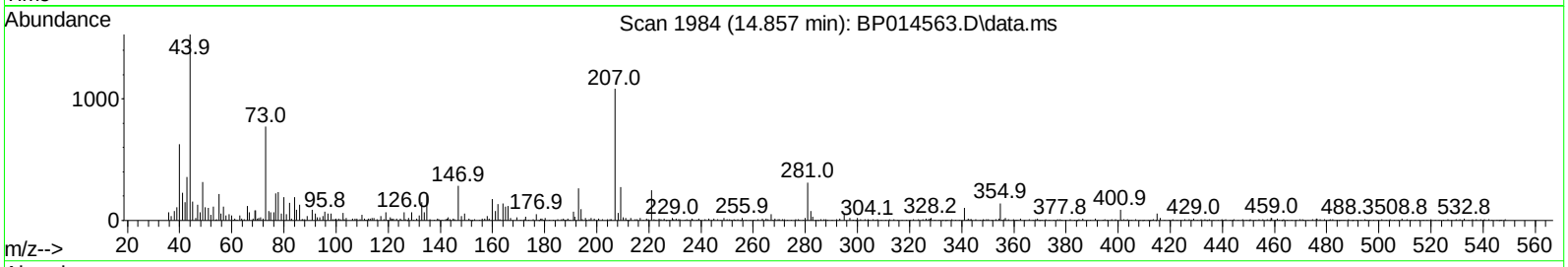

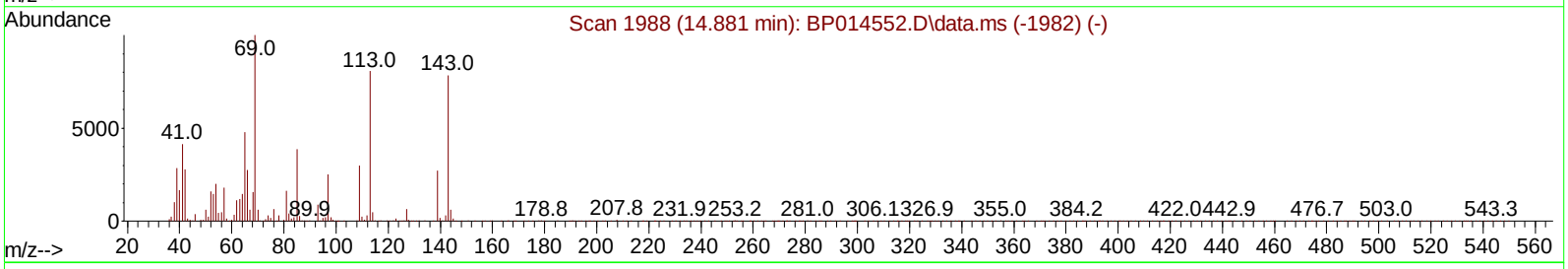

TIC: BP014563.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (-0.006) 0.01 ng/ul

response 27

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	40.54#
41.00	44.20	621.62#
42.00	31.50	405.41#

Quantitation Report (Qedit)

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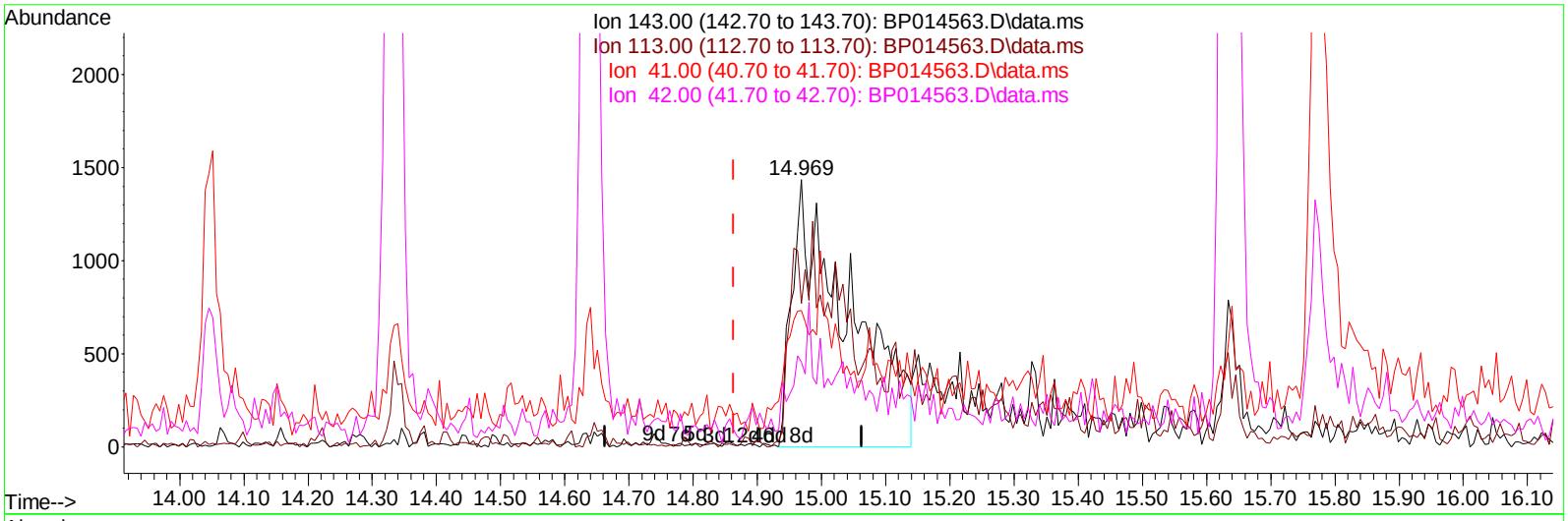

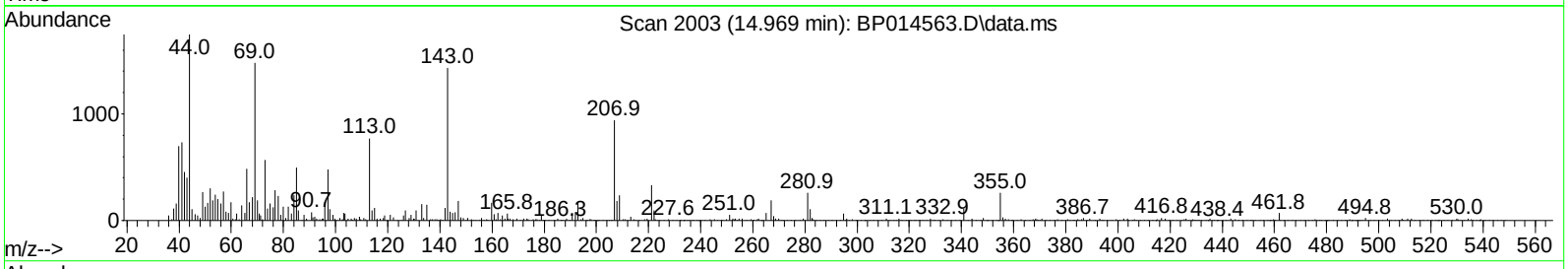

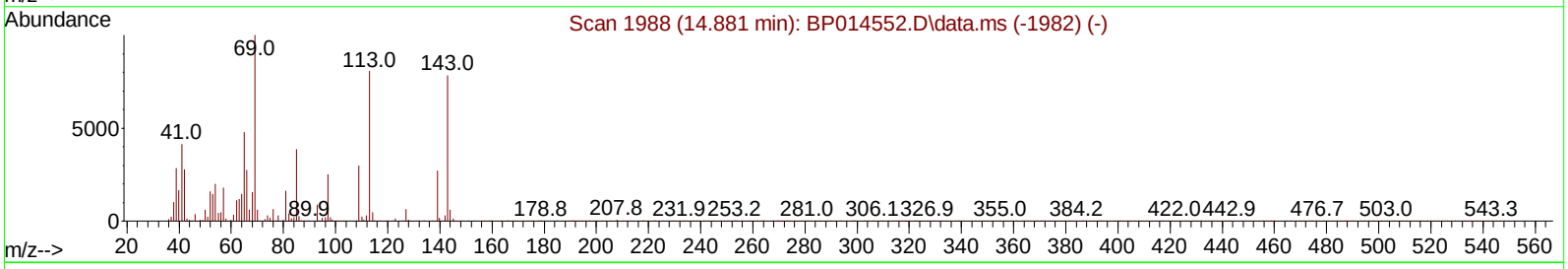

TIC: BP014563.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.969min (+ 0.106) 2.40 ng/ul m

response 8853

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	96.80	53.70#
41.00	44.20	51.05
42.00	31.50	31.87

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	107774	20.000	ng/ul	0.00
20) Naphthalene-d8	10.810	136	464760	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	301169	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	639993	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	578787	20.000	ng/ul	0.00
88) Perylene-d12	23.986	264	634202	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	12276	4.417	ng/uL	0.00
4) Pyridine-d5	3.817	84	48676m	6.835	ng/ul	0.02
7) Phenol-d5	7.169	99	35401	3.562	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.316	67	155925	24.450	ng/ul	0.00
11) 2-Chlorophenol-d4	7.522	132	118384	16.479	ng/ul	0.00
15) 4-Methylphenol-d8	8.722	113	73358	9.150	ng/ul	0.01
21) Nitrobenzene-d5	9.169	128	84968	22.646	ng/ul	0.00
24) 2-Nitrophenol-d4	9.893	143	79345	18.469	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	135241	17.285	ng/ul	0.01
31) 4-Chloroaniline-d4	10.969	131	173093	16.795	ng/ul	0.01
46) Dimethylphthalate-d6	14.051	166	653031	26.742	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	683324	25.219	ng/ul	0.00
54) 4-Nitrophenol-d4	14.969	143	8853m	2.400	ng/ul	0.11
60) Fluorene-d10	15.634	176	534611	26.253	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	66319	14.083	ng/ul	0.01
73) Anthracene-d10	17.504	188	877962	28.346	ng/ul	0.00
81) Pyrene-d10	19.728	212	1048514	26.970	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	1021185	30.307	ng/ul	0.00

Target Compounds Qvalue

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

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