Quantitation Report (Qedit)

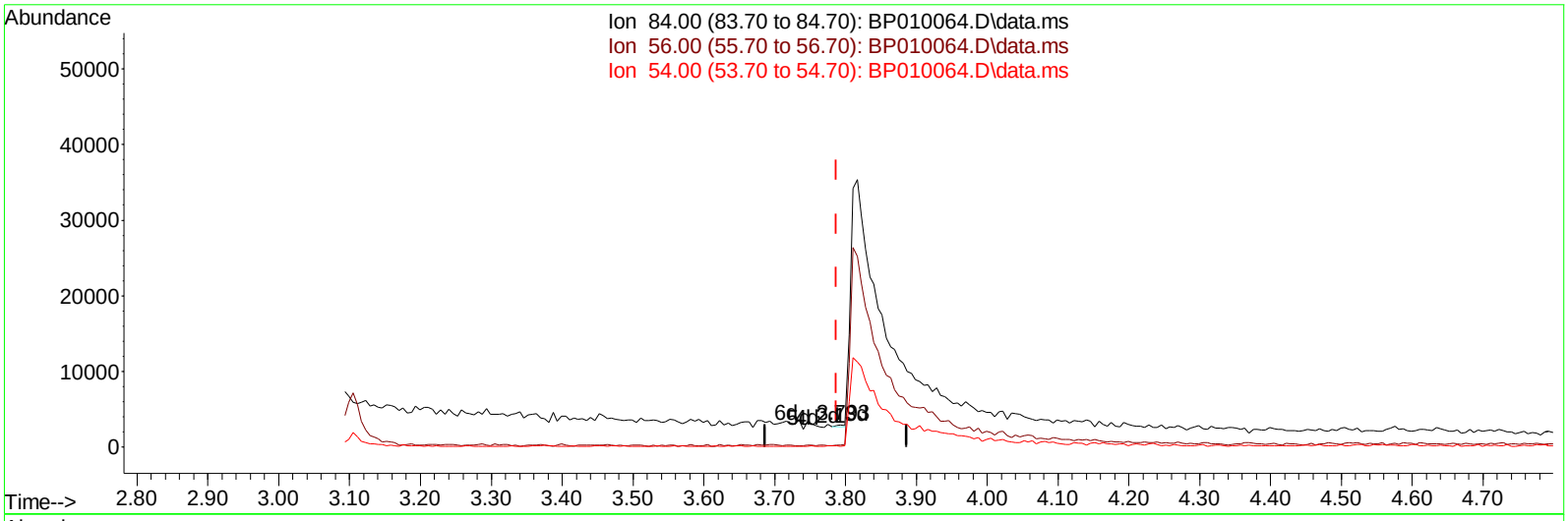
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010064.D
Acq On : 22 Apr 2022 20:34
Operator : CG/JU
Sample : N2490-10
Misc :
ALS Vial : 60 Sample Multiplier: 1

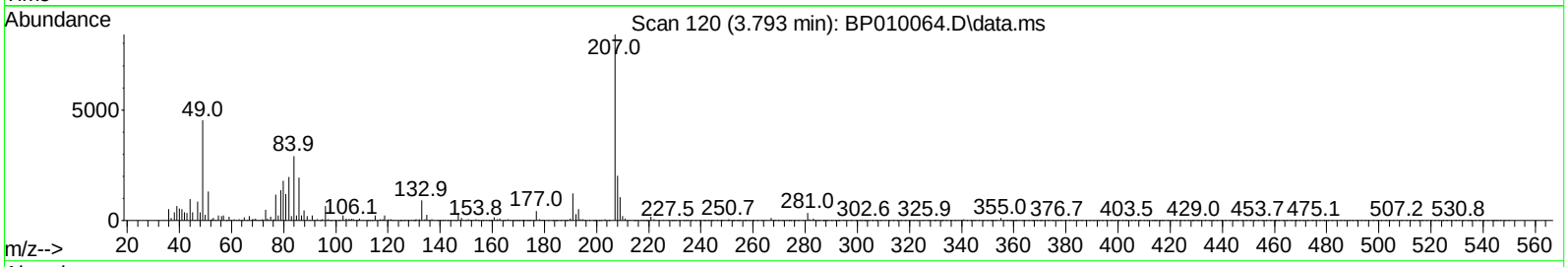
Quant Time: Apr 22 23:31:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

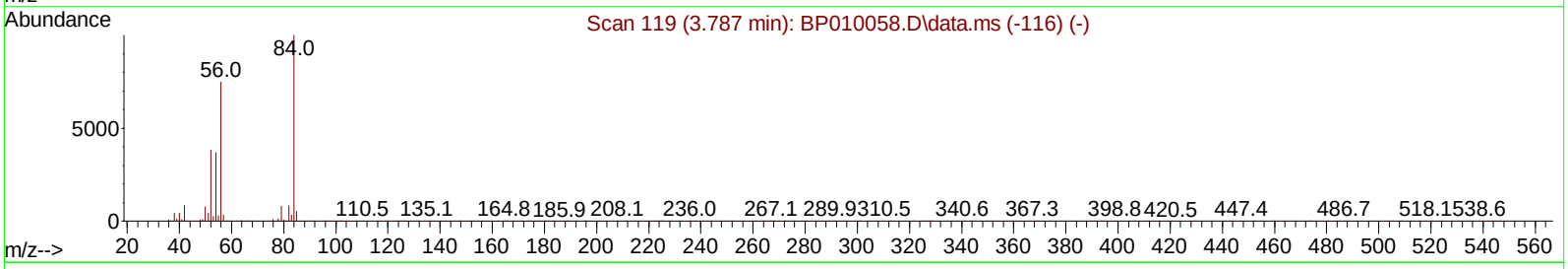
Instrument :
BNA_P
ClientSampleId :
C0J27

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/25/2022
Supervised By :mohammad ahmed 04/26/2022







TIC: BP010064.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.006) 0.03 ng/ul

response 262

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	7.22#
54.00	24.80	0.62#
0.00	0.00	0.00

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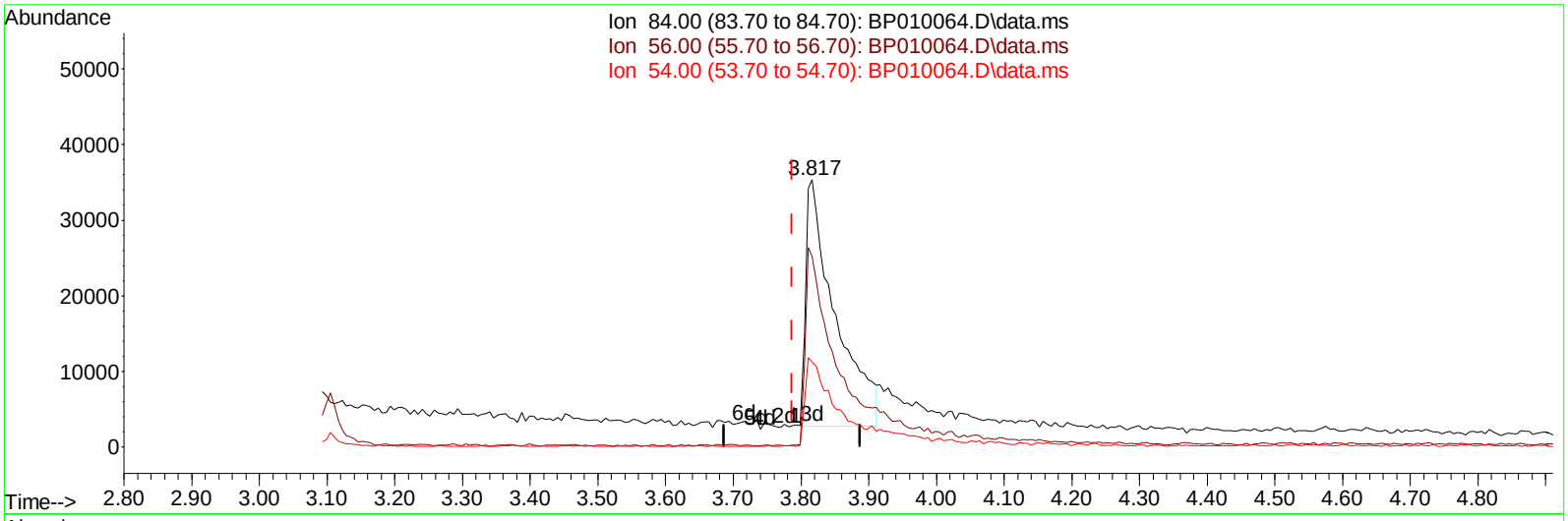

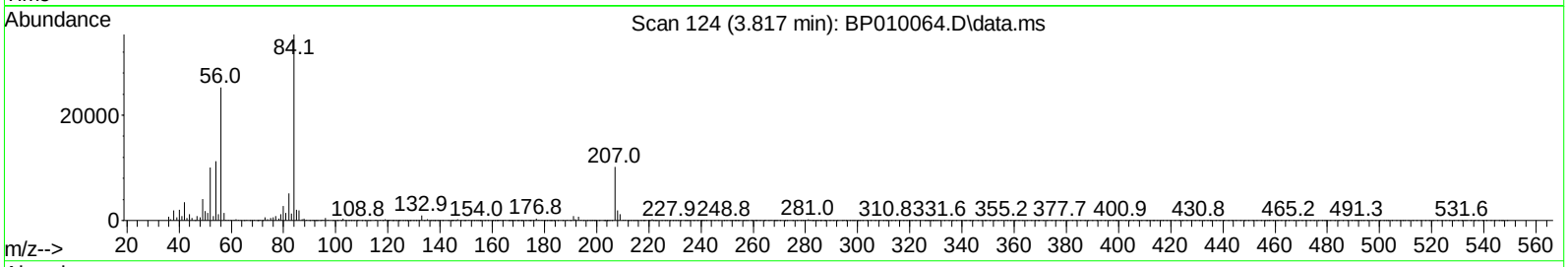

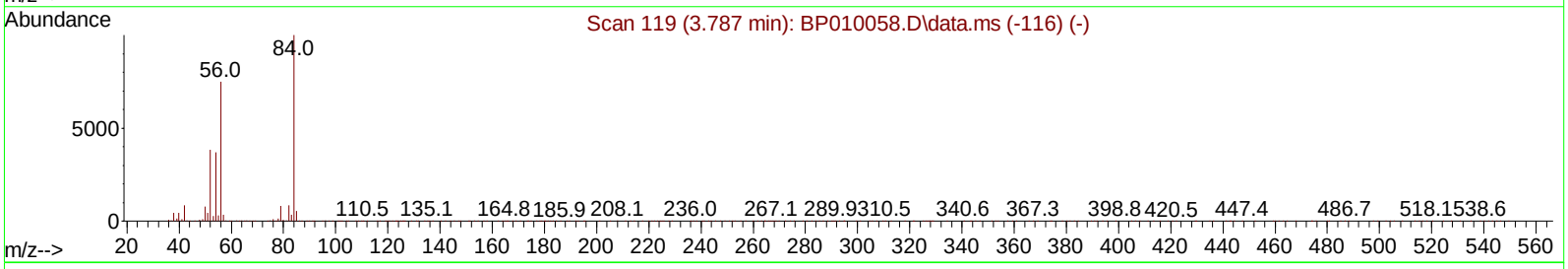

TIC: BP010064.D\data.ms

(4) Pyridine-d5 (S)

3.817min (+ 0.030) 11.55 ng/ul m

response 97828

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	30.50	71.44#
54.00	24.80	31.80#
0.00	0.00	0.00

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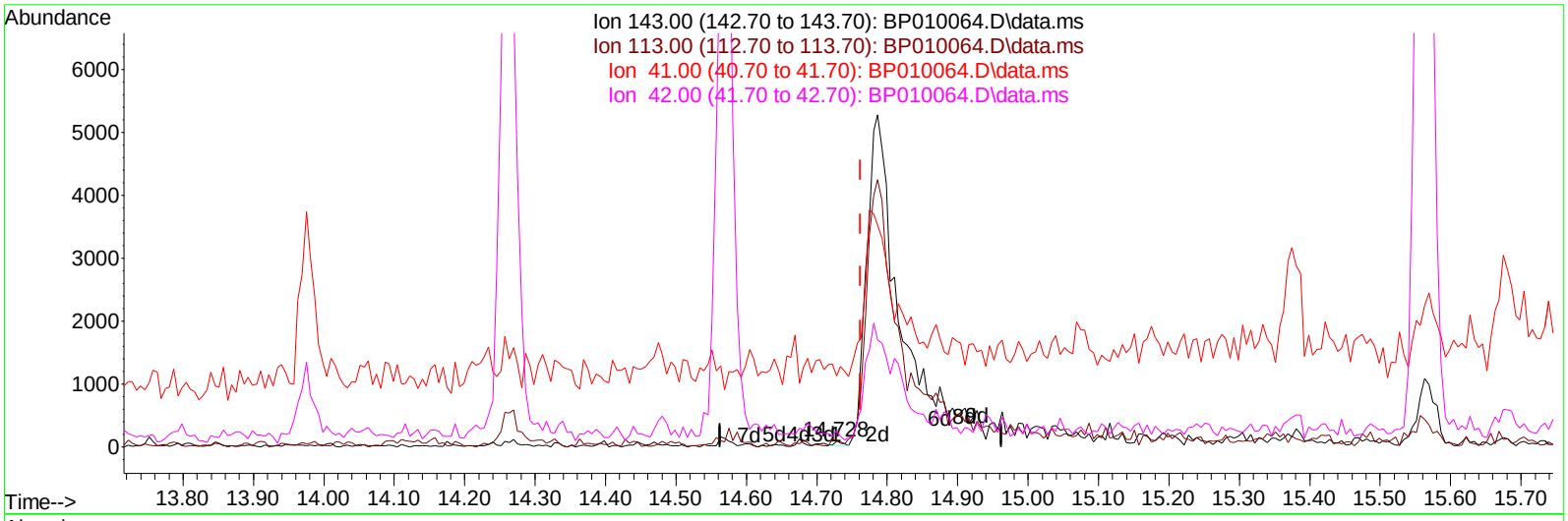

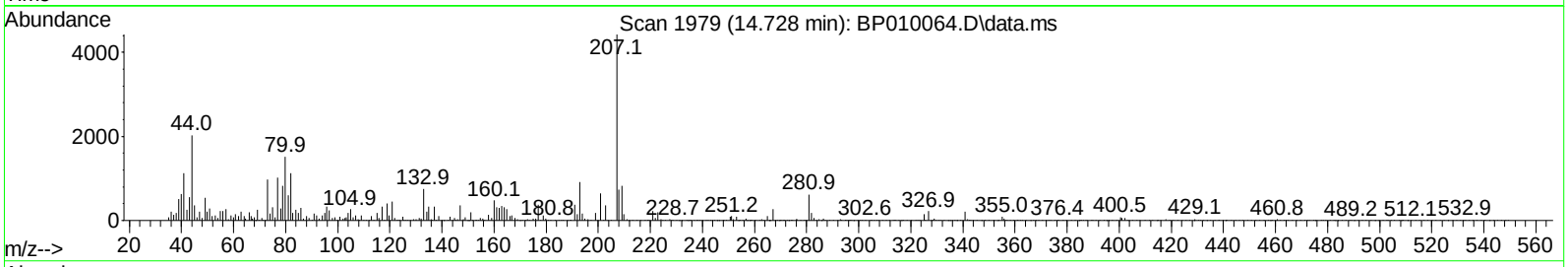

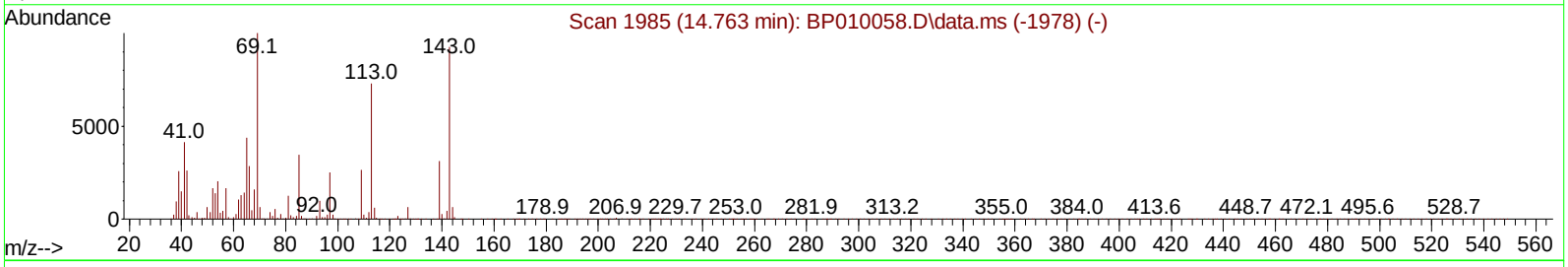

TIC: BP010064.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.035) 0.01 ng/ul

response 62

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	126.74
41.00	23.20	1308.14#
42.00	18.00	293.02#

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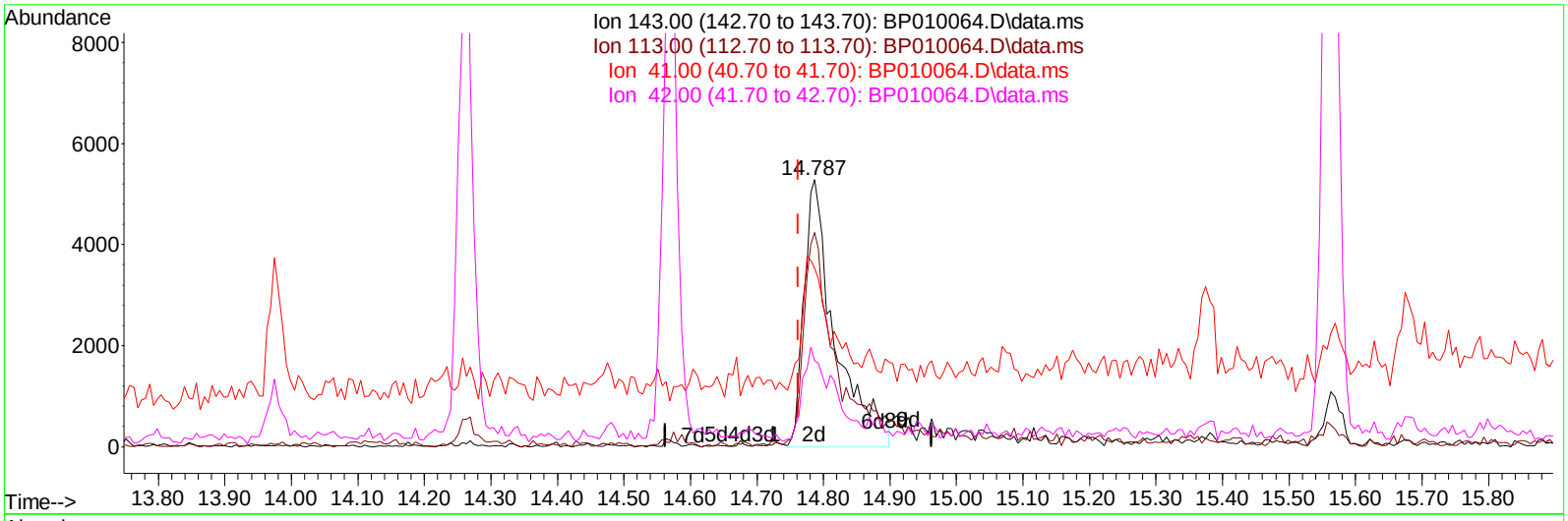

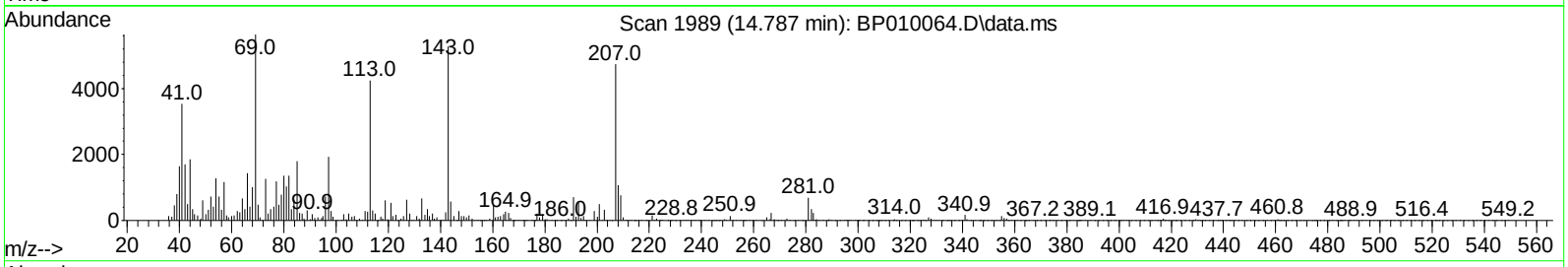

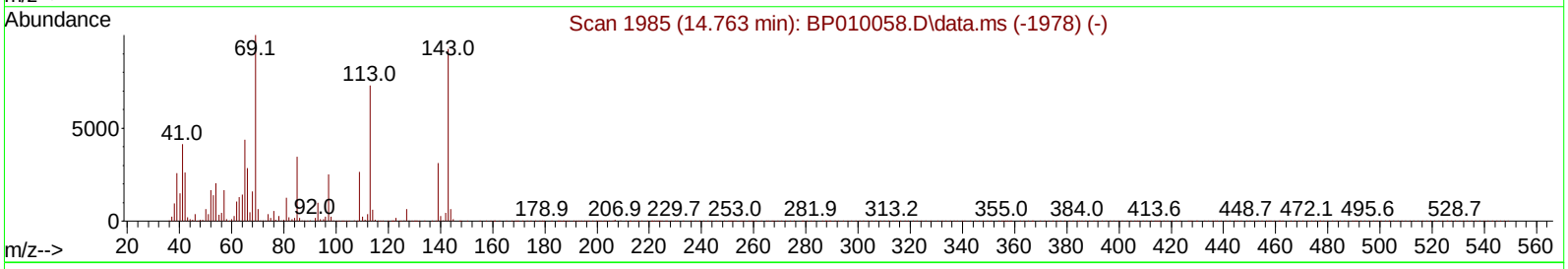

TIC: BP010064.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.787min (+ 0.024) 2.49 ng/ul m

response 17260

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.39#
41.00	23.20	66.81#
42.00	18.00	32.20#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	131326	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	654969	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	483096	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1029100	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	712146	20.000	ng/ul	# 0.00
88) Perylene-d12	23.863	264	601551	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	16497	5.478	ng/uL	0.00
4) Pyridine-d5	3.817	84	97828m	11.547	ng/ul	0.03
7) Phenol-d5	7.105	99	76008	6.363	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	213779	29.825	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	212935	24.070	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	145270	14.678	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	144926	27.821	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	100945	17.743	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	273559	24.699	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	315763	19.884	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1106591	29.447	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1263592	27.783	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	17260m	2.495	ng/ul	0.02
60) Fluorene-d10	15.563	176	988181	30.569	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	21014	3.229	ng/ul	0.00
73) Anthracene-d10	17.422	188	1504310	30.215	ng/ul	0.00
81) Pyrene-d10	19.651	212	1623179	40.078	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	958935	30.407	ng/ul	0.00

Target Compounds Qvalue

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

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