Quantitation Report (Qedit)

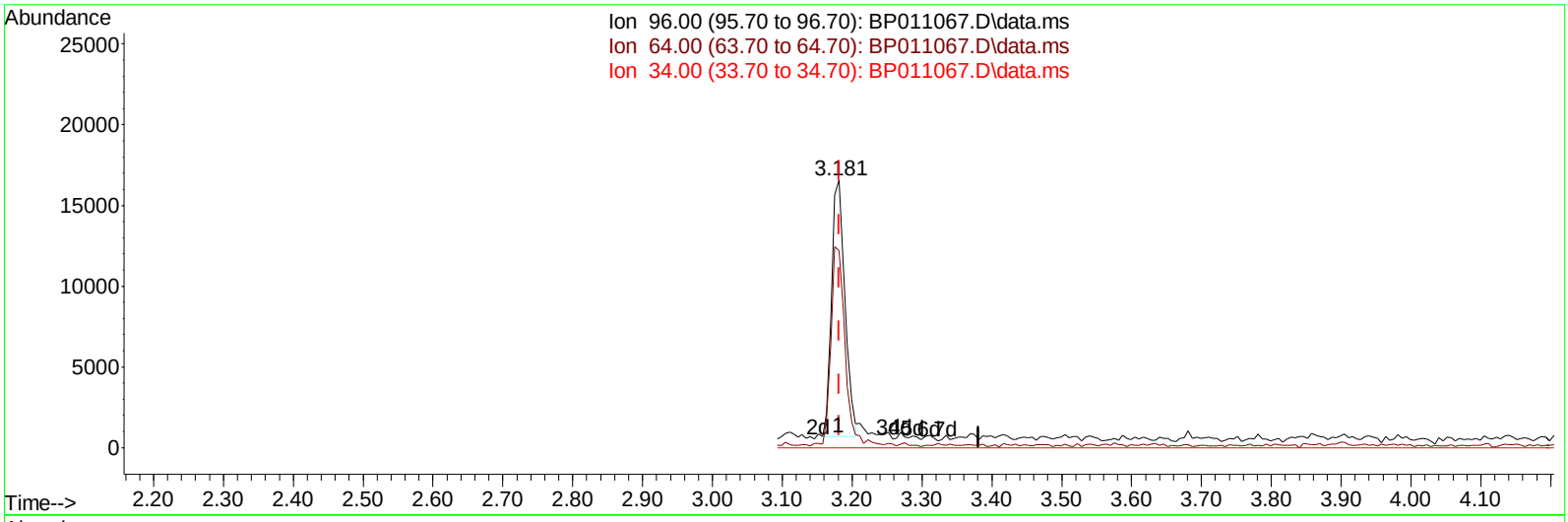
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
Data File : BP011067.D
Acq On : 21 Jul 2022 17:54
Operator : CG/JU
Sample : N3709-12
Misc :
ALS Vial : 10 Sample Multiplier: 1

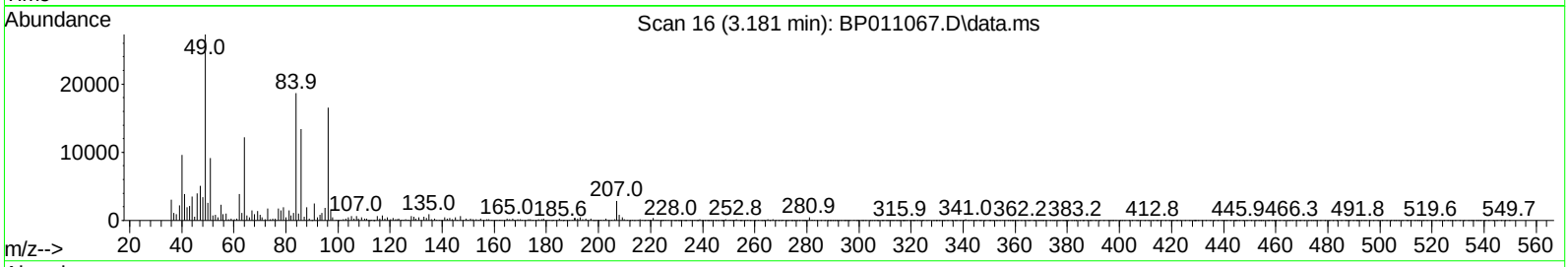
Quant Time: Jul 21 23:24:57 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 19 03:25:54 2022
Response via : Initial Calibration

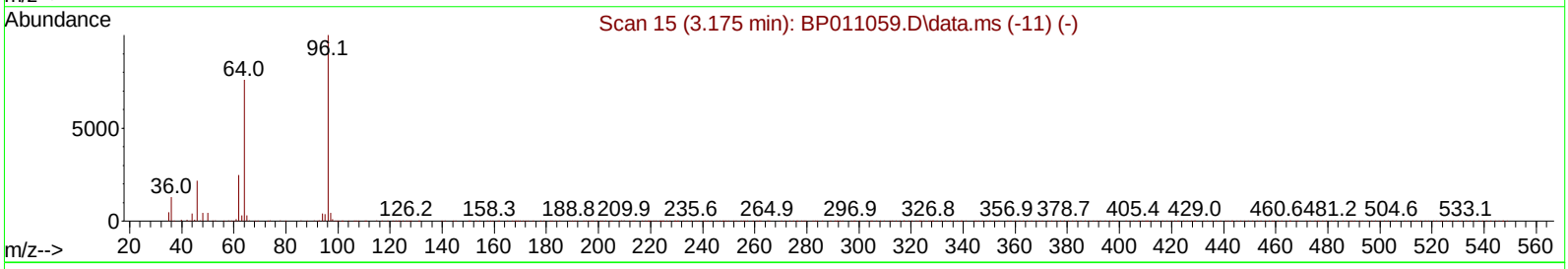
Instrument :
BNA_P
ClientSampleId :
H0B26

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/22/2022
Supervised By :mohammad ahmed 07/22/2022







TIC: BP011067.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.181min (-0.000) 2.00 ng/uL

response 20747

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	73.60
34.00	0.00	0.00
0.00	0.00	0.00

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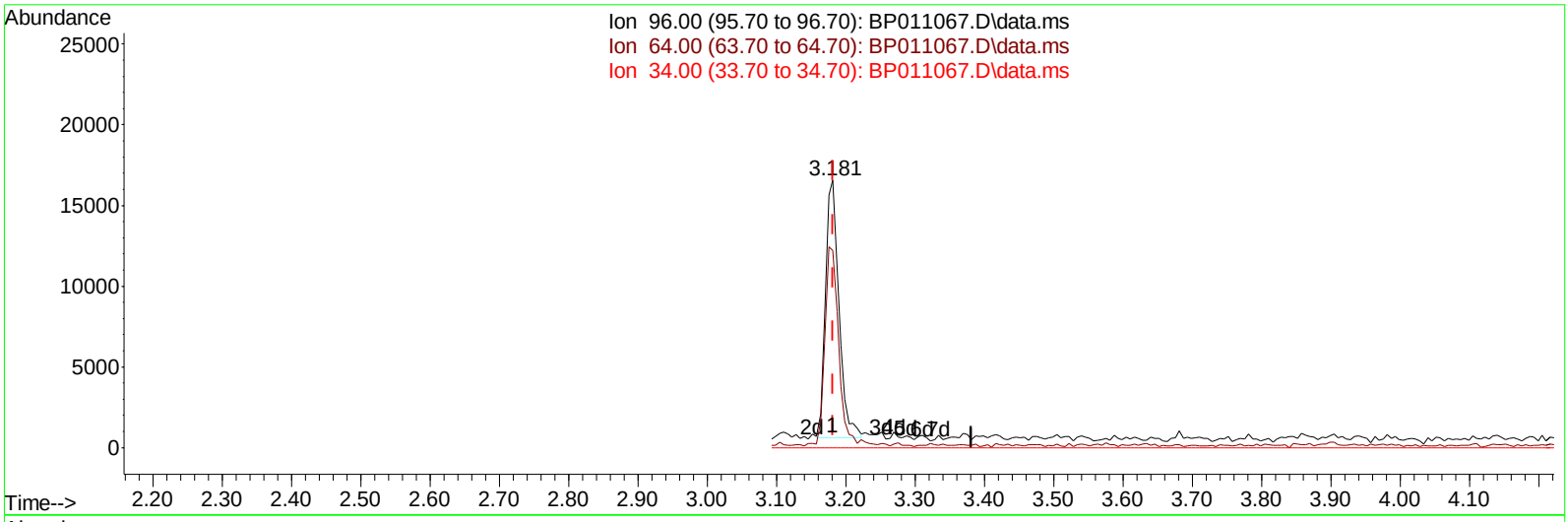

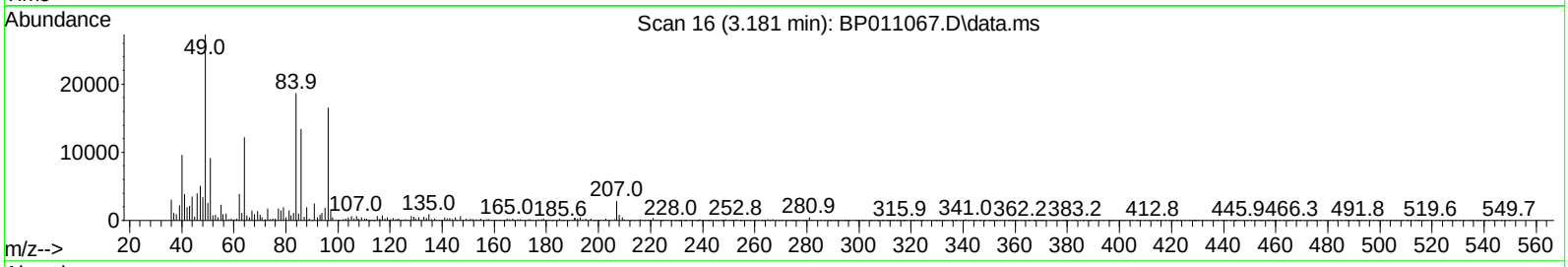

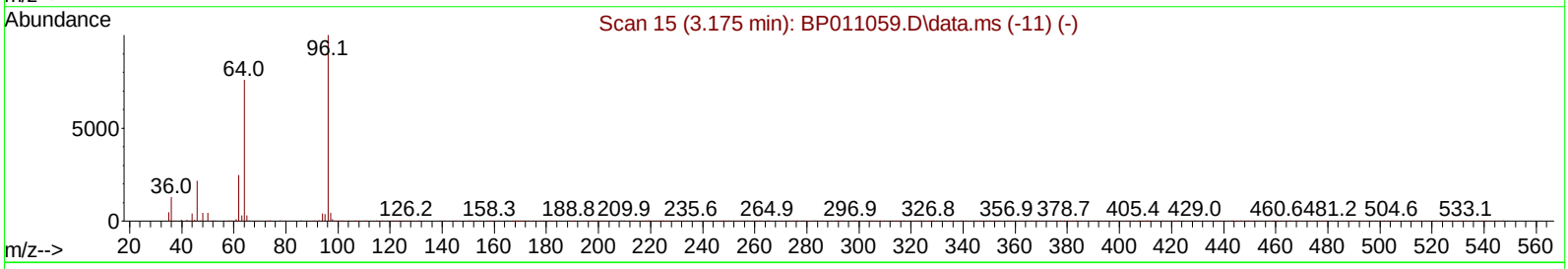

TIC: BP011067.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.181min (-0.000) 2.08 ng/uL m

response 21598

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	73.60
34.00	0.00	0.00
0.00	0.00	0.00

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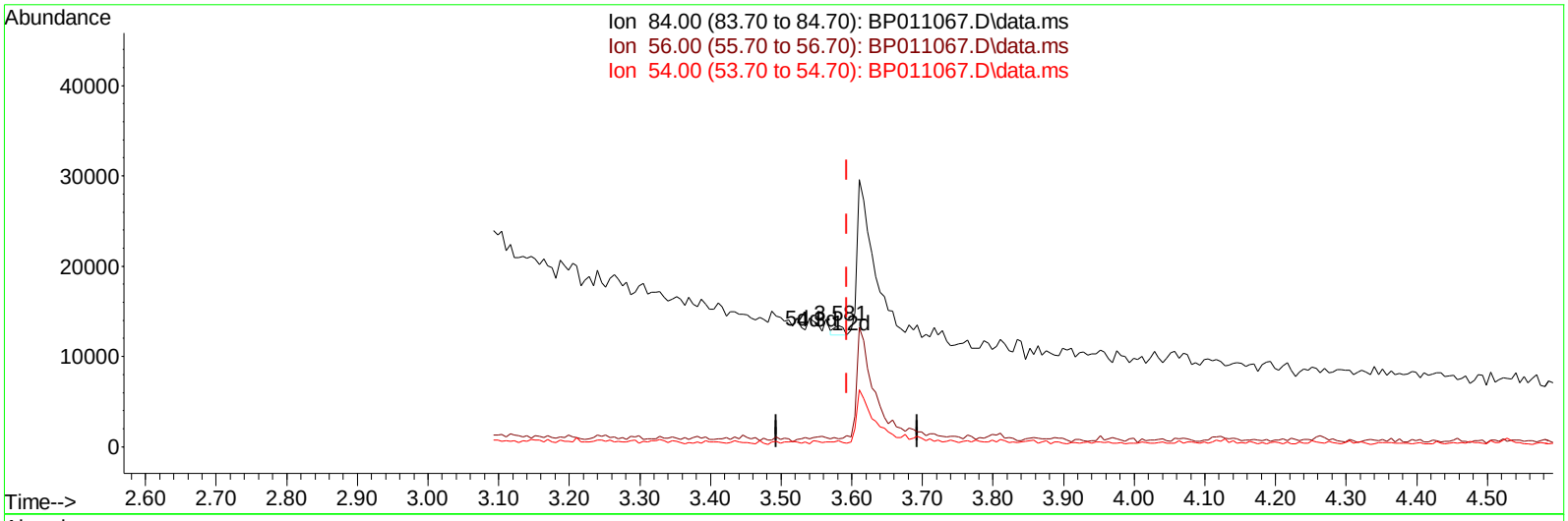

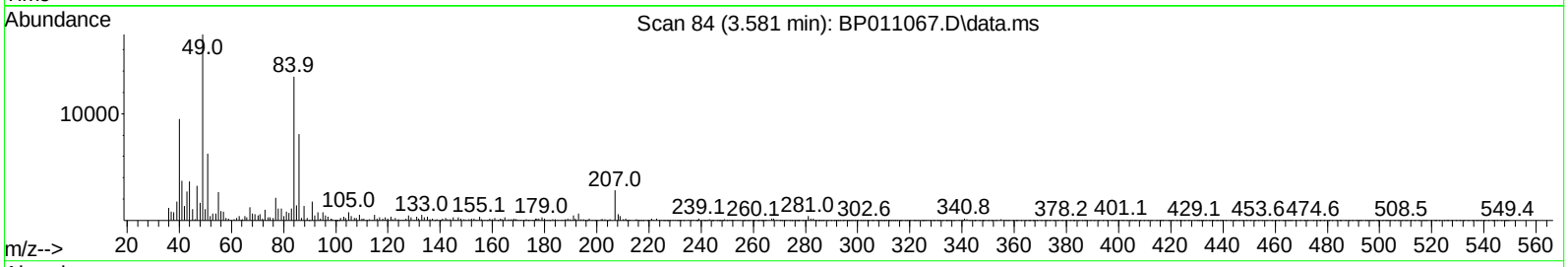

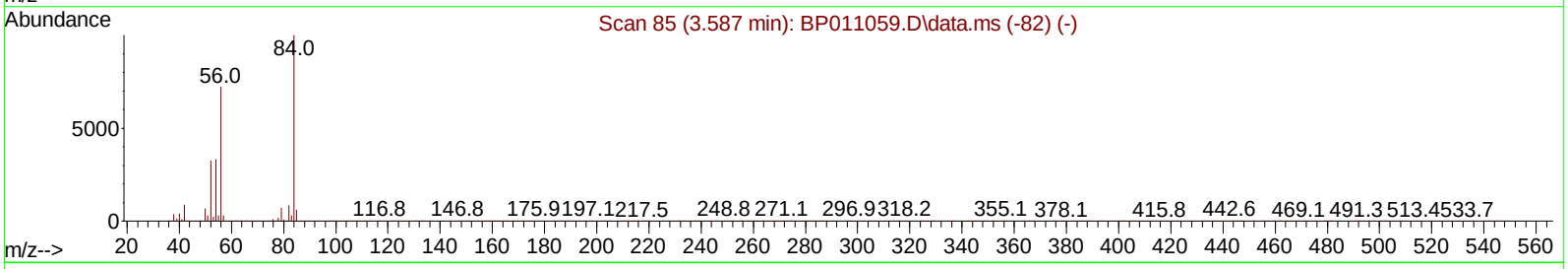

TIC: BP011067.D\data.ms

(4) Pyridine-d5 (S)

3.581min (-0.012) 0.04 ng/ul

response 946

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	6.60#
54.00	44.80	4.73#
0.00	0.00	0.00

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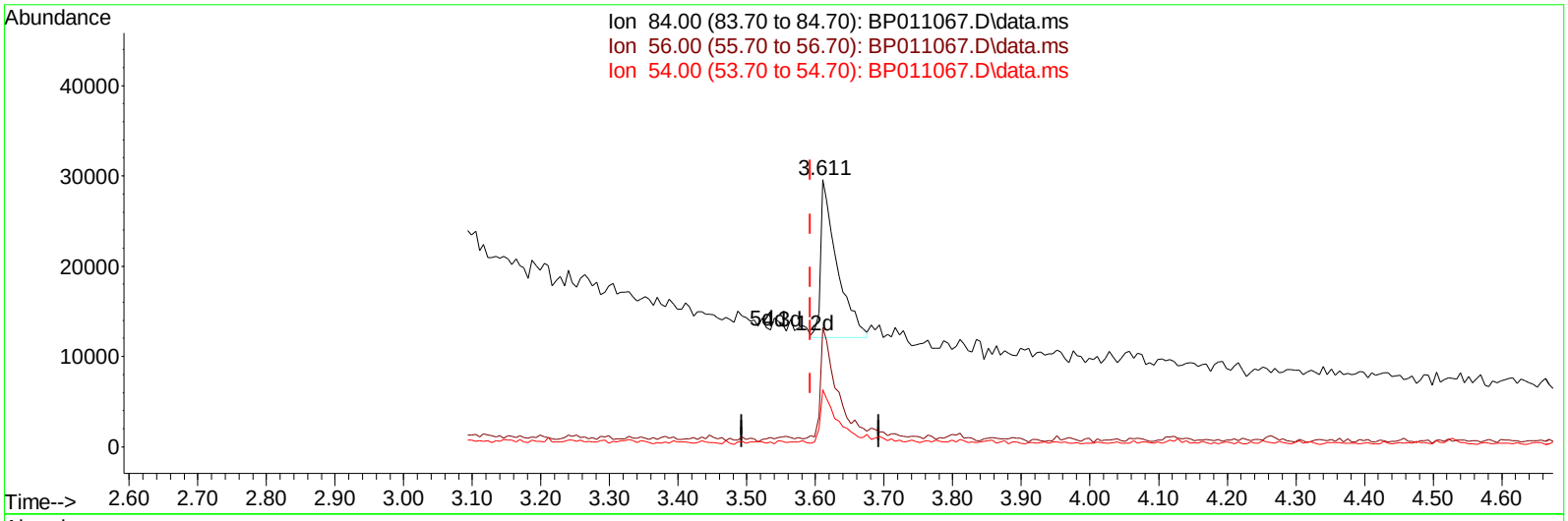

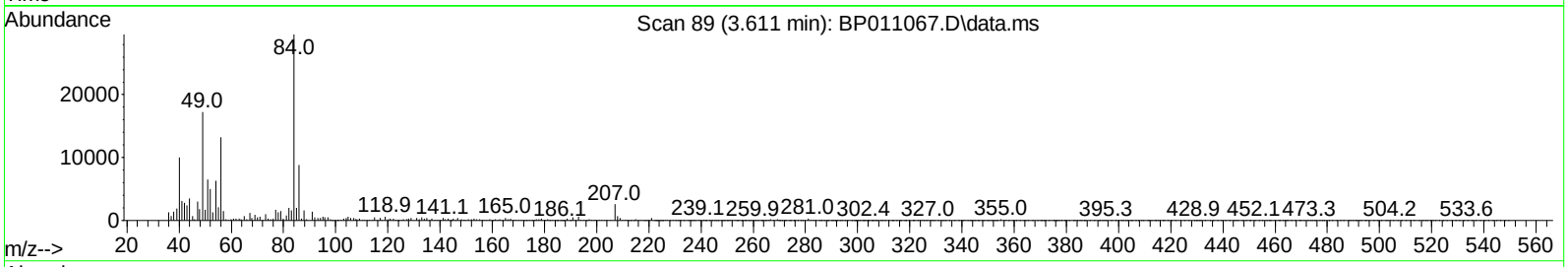

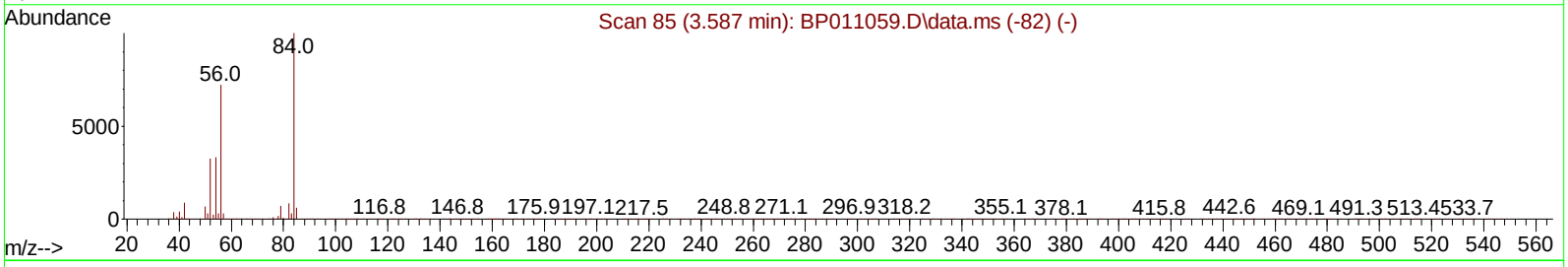

TIC: BP011067.D\data.ms

(4) Pyridine-d5 (S)

3.611min (+ 0.018) 1.08 ng/ul m

response 28976

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	44.66#
54.00	44.80	21.31#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	374514	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	1545871	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.334	164	875840	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1717682	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1585231	20.000	ng/ul	# 0.00
88) Perylene-d12	23.574	264	1616813	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	21598m	2.079	ng/uL	0.00
4) Pyridine-d5	3.611	84	28976m	1.079	ng/ul	0.02
7) Phenol-d5	6.852	99	298872	9.722	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	258688	13.086	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	214219	8.624	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	204298	8.301	ng/ul	-0.01
21) Nitrobenzene-d5	8.834	128	156217	13.867	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	30474	2.734	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	98071	4.811	ng/ul	0.00
31) 4-Chloroaniline-d4	10.610	131	370499	11.024	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	828373	13.838	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	1062060	15.037	ng/ul	0.00
54) 4-Nitrophenol-d4	14.545	143	37	0.003	ng/ul	-0.01
60) Fluorene-d10	15.334	176	747247	14.686	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	78	0.009	ng/ul	0.00
73) Anthracene-d10	17.198	188	1049075	14.366	ng/ul	0.00
81) Pyrene-d10	19.457	212	1276636	15.334	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.421	264	1171659	15.088	ng/ul	0.00
Target Compounds						Qvalue

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

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