Quantitation Report (Qedit)

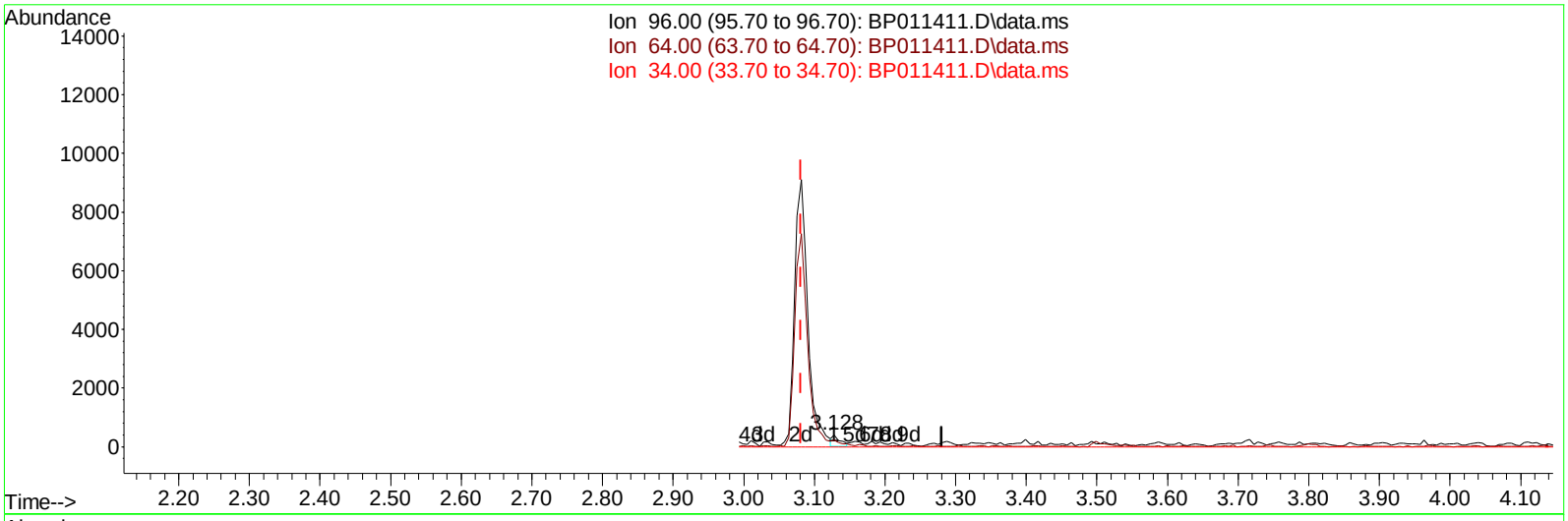
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011411.D
Acq On : 12 Aug 2022 05:25
Operator : CG/JU
Sample : N4129-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

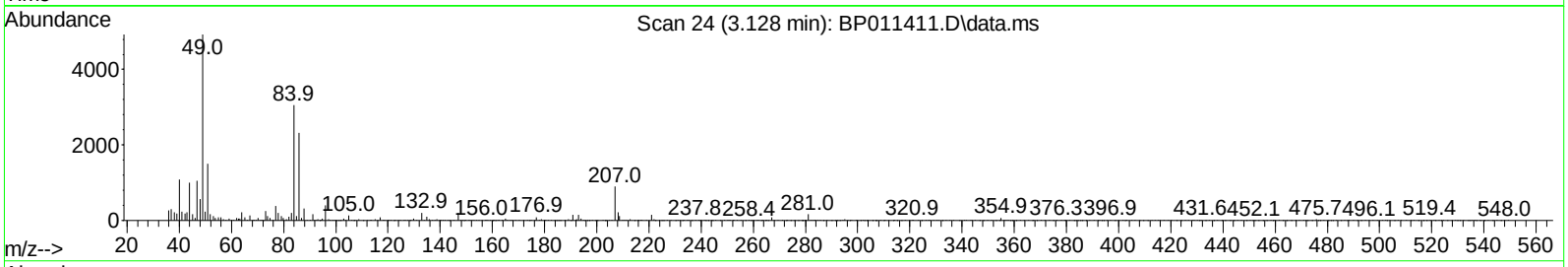
Quant Time: Aug 12 07:16:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

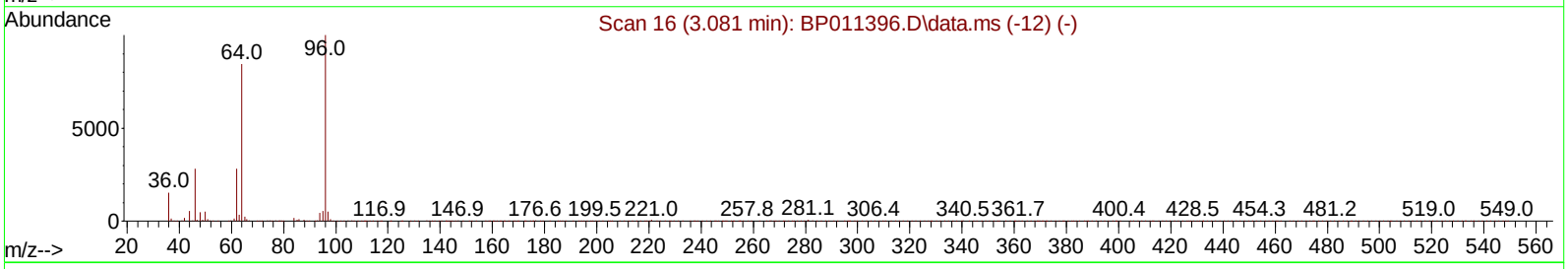
Instrument :
BNA_P
ClientSampleId :
BGBW5

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011411.D\data.ms

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

3.128min (+ 0.047) 0.11 ng/uL

response 296

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	54.21
34.00	0.00	0.00
0.00	0.00	0.00

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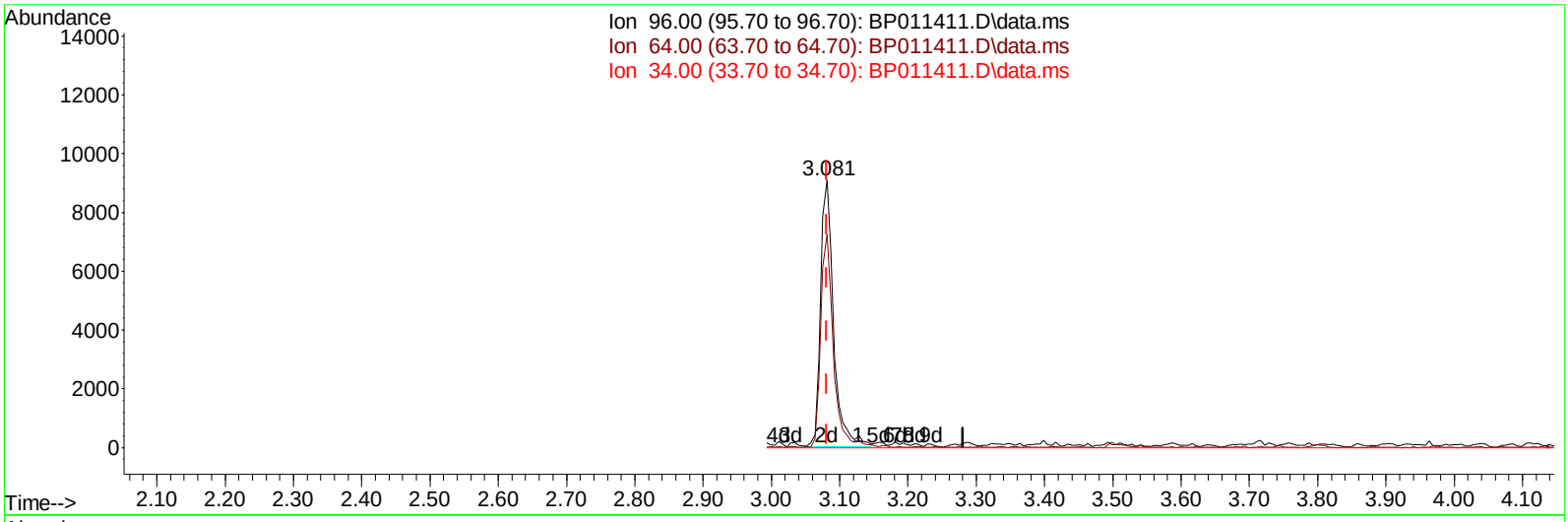

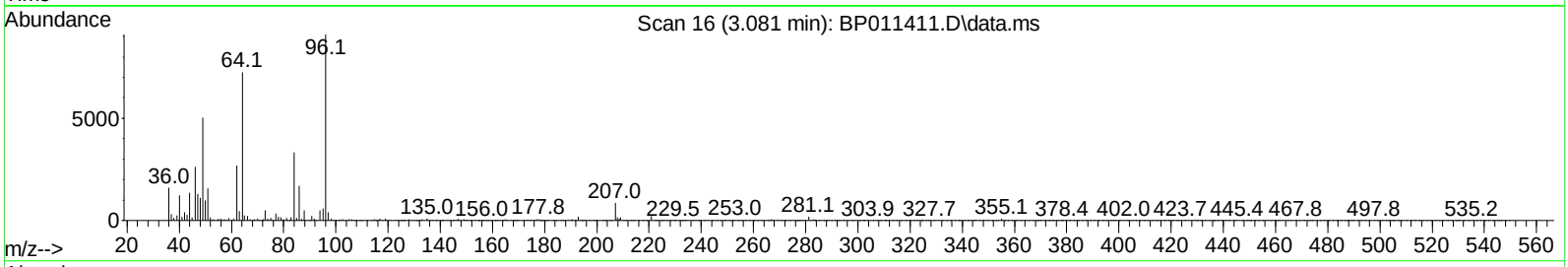

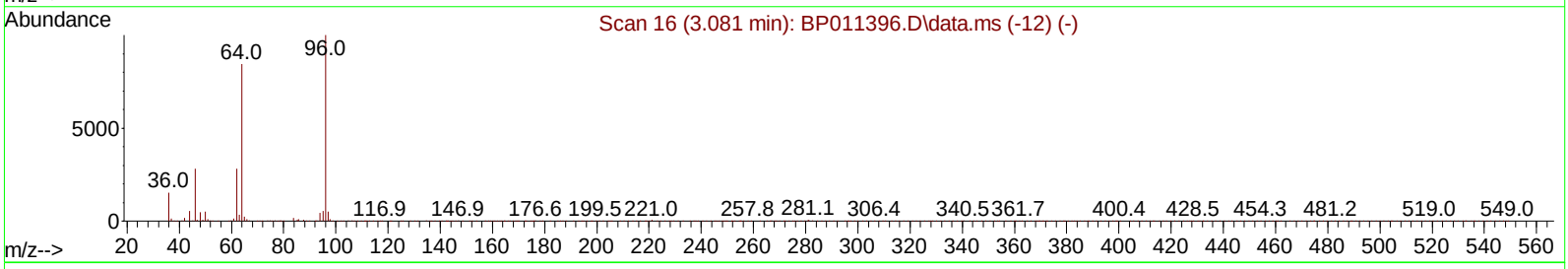

TIC: BP011411.D\data.ms

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

3.081min (0.000) 4.65 ng/uL m

response 12034

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	79.80#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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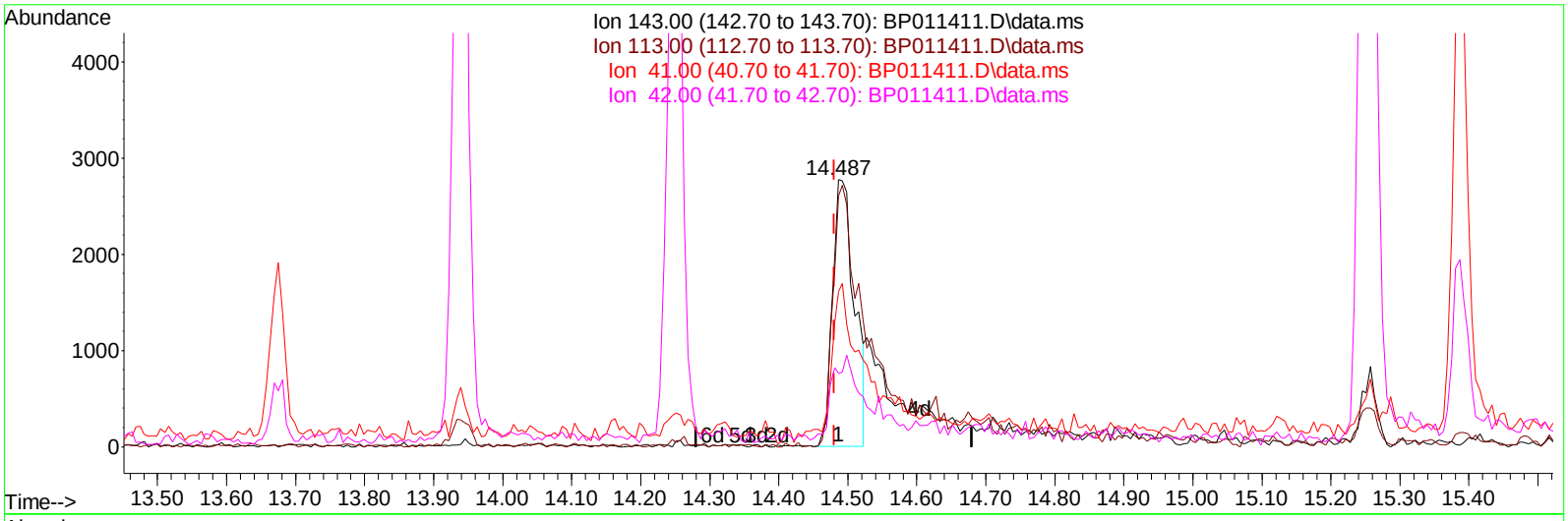

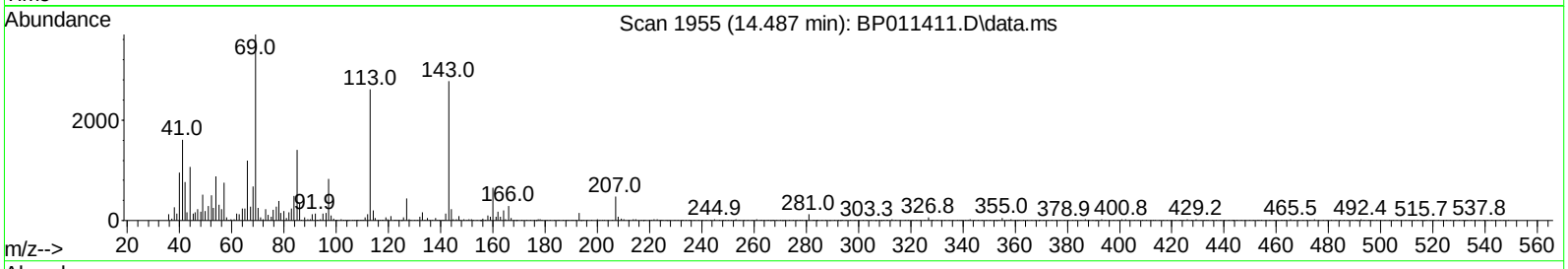

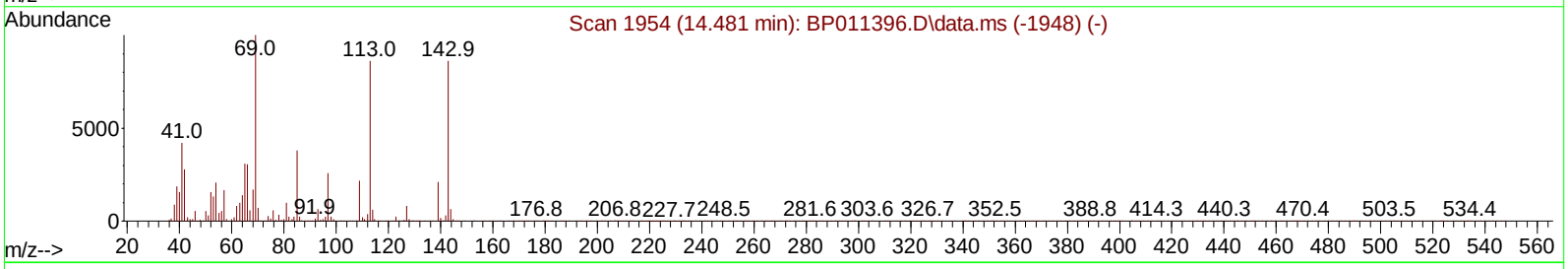

TIC: BP011411.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.006) 2.48 ng/ul

response 6086

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.05#
41.00	54.40	58.18
42.00	35.50	27.58#

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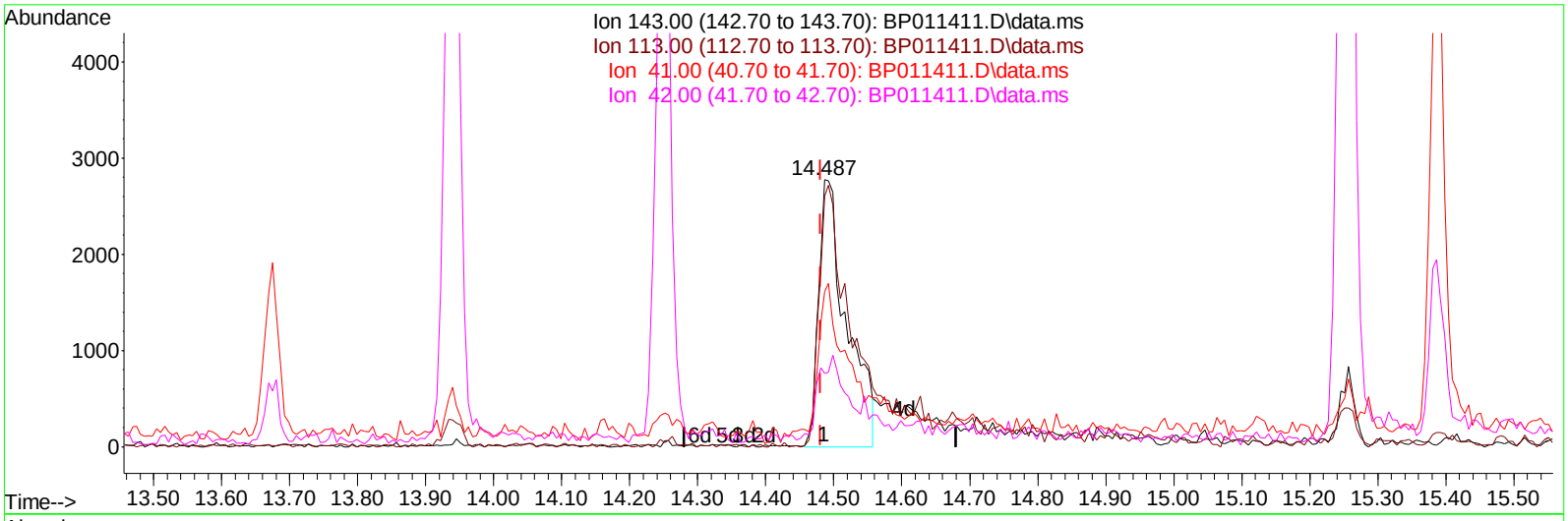

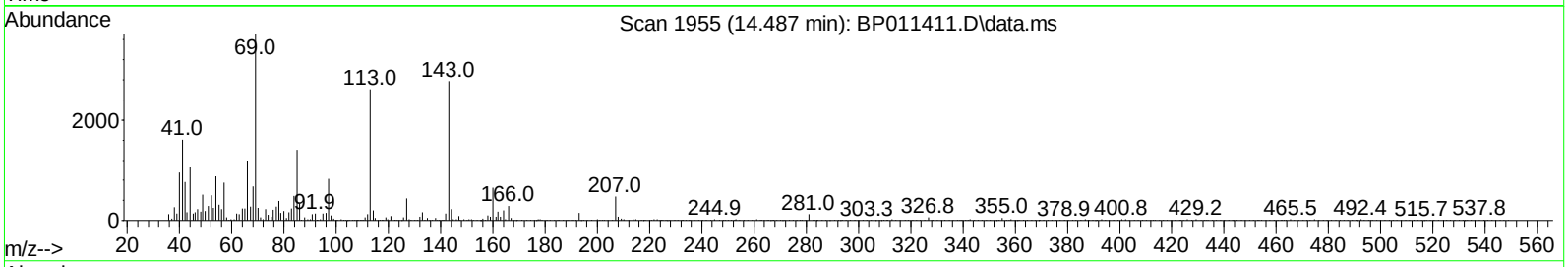

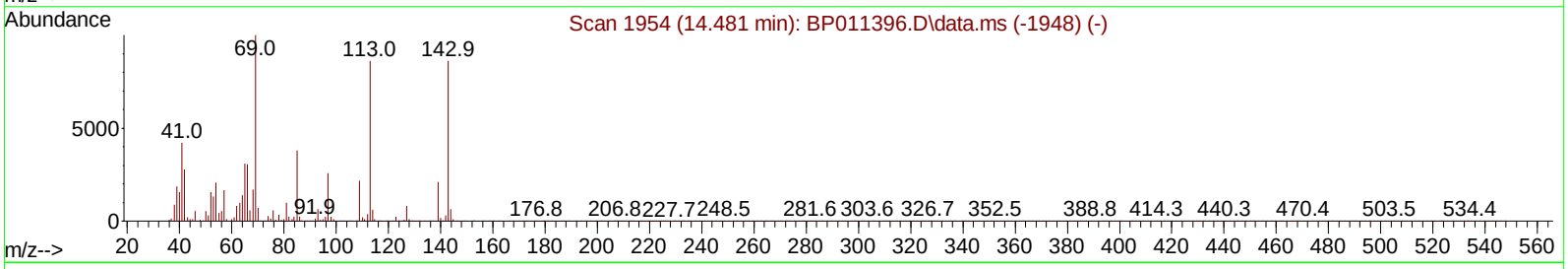

TIC: BP011411.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.487min (+ 0.006) 3.23 ng/ul m

response 7920

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	94.05#
41.00	54.40	58.18
42.00	35.50	27.58#

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Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.575	152	79277	20.000	ng/ul		0.00
20) Naphthalene-d8	10.363	136	338811	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.251	164	193738	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.022	188	405432	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.151	240	391278	20.000	ng/ul		0.00
88) Perylene-d12	23.457	264	413206	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.081	96	12034m	4.653	ng/uL		0.00
4) Pyridine-d5	3.499	84	51934	7.961	ng/ul		0.01
7) Phenol-d5	6.758	99	43921	6.260	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	156191	31.398	ng/ul		0.00
11) 2-Chlorophenol-d4	7.110	132	139673	25.373	ng/ul		0.00
15) 4-Methylphenol-d8	8.299	113	86197	15.647	ng/ul		0.00
21) Nitrobenzene-d5	8.740	128	86242	37.401	ng/ul		0.00
24) 2-Nitrophenol-d4	9.457	143	83817	39.703	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	9.993	165	126652	29.707	ng/ul		0.00
31) 4-Chloroaniline-d4	10.516	131	176560	23.688	ng/ul		0.00
46) Dimethylphthalate-d6	13.675	166	528218	39.295	ng/ul		0.00
49) Acenaphthylene-d8	13.939	160	598049	36.959	ng/ul		0.00
54) 4-Nitrophenol-d4	14.487	143	7920m	3.226	ng/ul		0.00
60) Fluorene-d10	15.257	176	436185	37.942	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	62390	33.130	ng/ul		0.00
73) Anthracene-d10	17.122	188	748035	41.048	ng/ul		0.00
81) Pyrene-d10	19.386	212	872974	40.496	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.310	264	855055	41.033	ng/ul		0.00

Target Compounds Qvalue

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

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