Quantitation Report (Qedit)

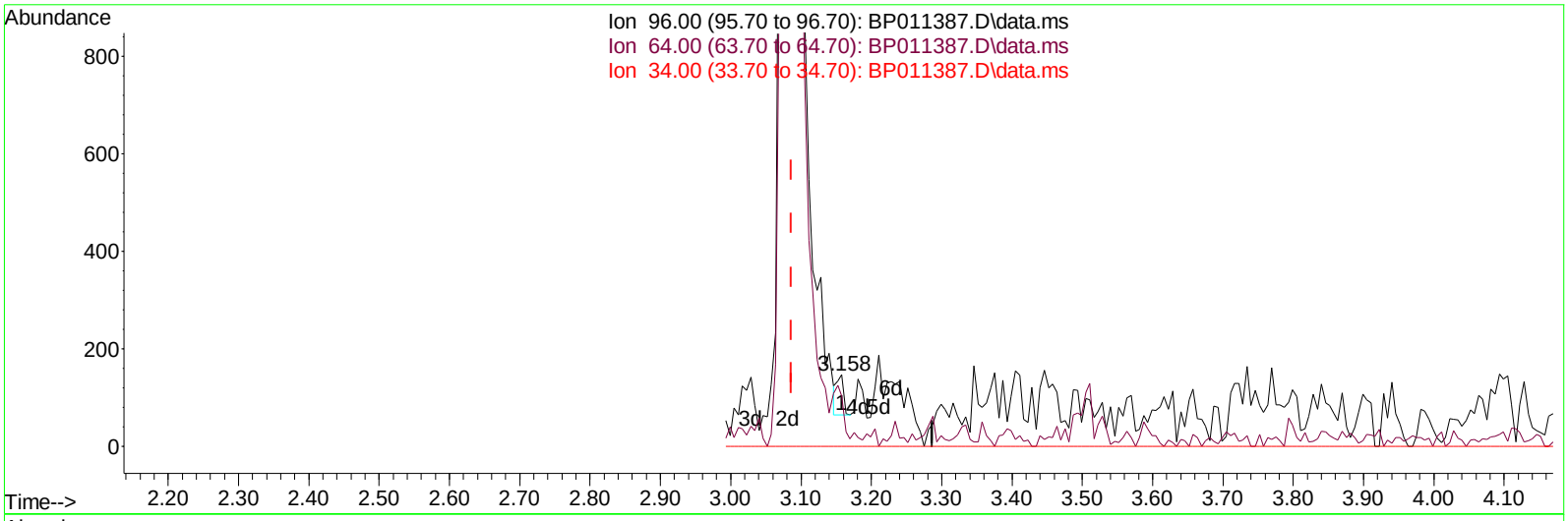
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011387.D
Acq On : 11 Aug 2022 15:02
Operator : CG/JU
Sample : N4128-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

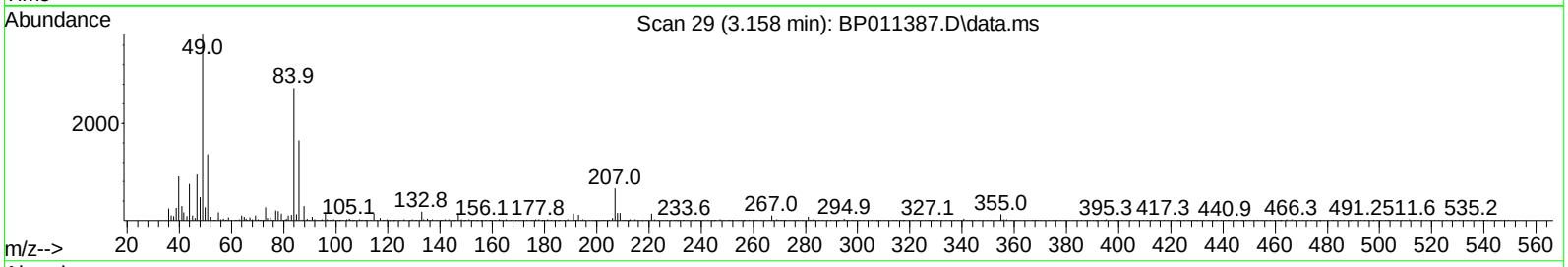
Quant Time: Aug 11 16:17:15 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

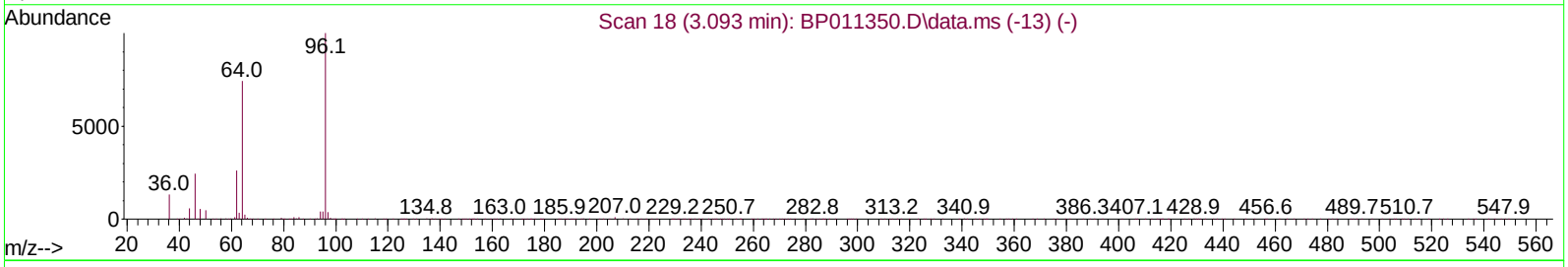
Instrument :
BNA_P
ClientSampleId :
BGBZ9

Manual IntegrationsAPPROVED

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







TIC: BP011387.D\data.ms

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

3.158min (+ 0.071) 0.02 ng/uL

response 55

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	61.18
34.00	0.00	0.00
0.00	0.00	0.00

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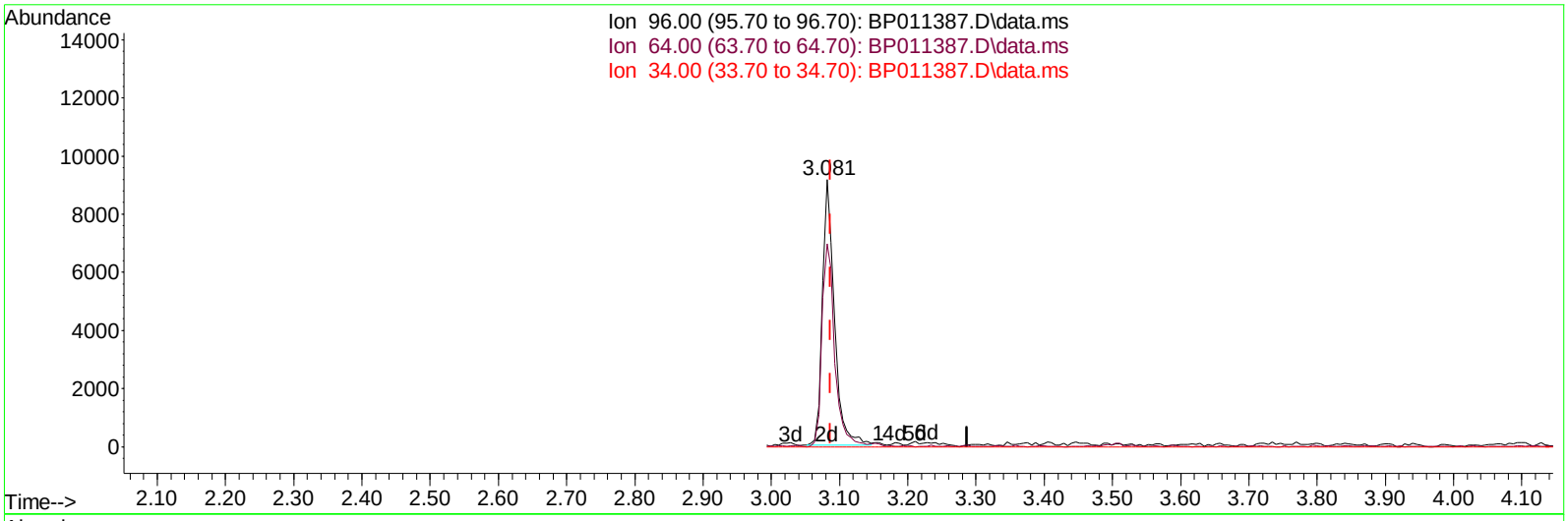

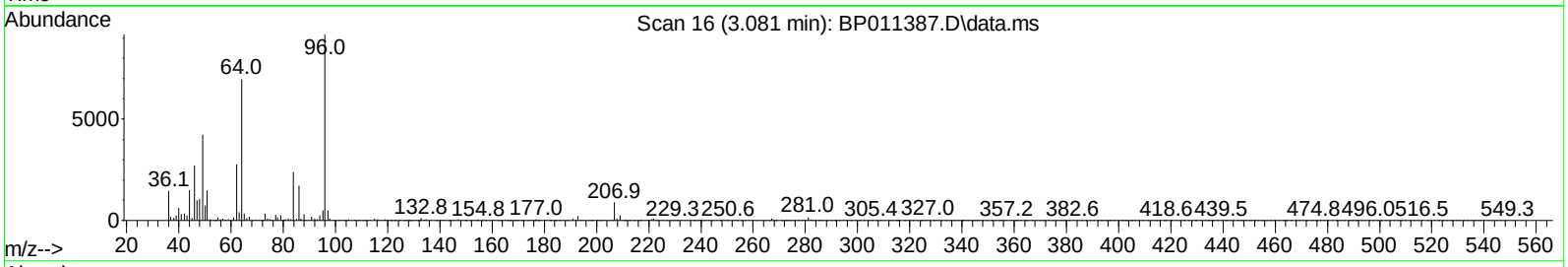

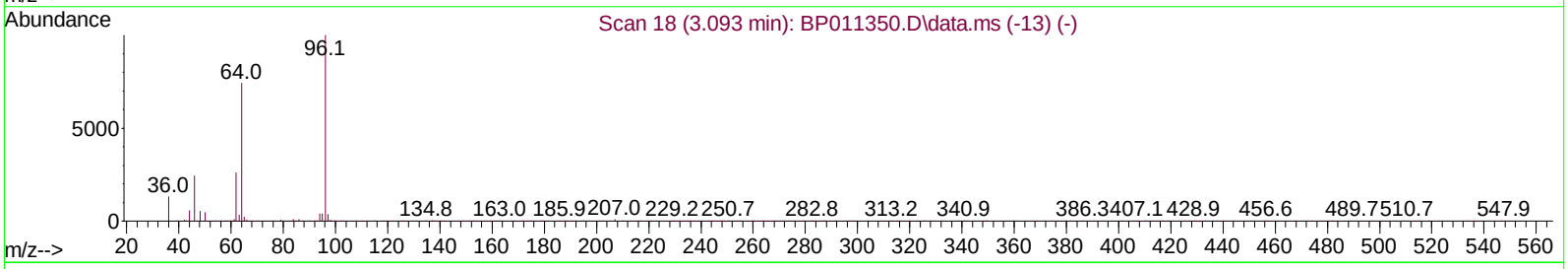

TIC: BP011387.D\data.ms

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

3.081min (-0.006) 4.07 ng/uL m

response 11189

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	76.06#
34.00	0.00	0.00
0.00	0.00	0.00

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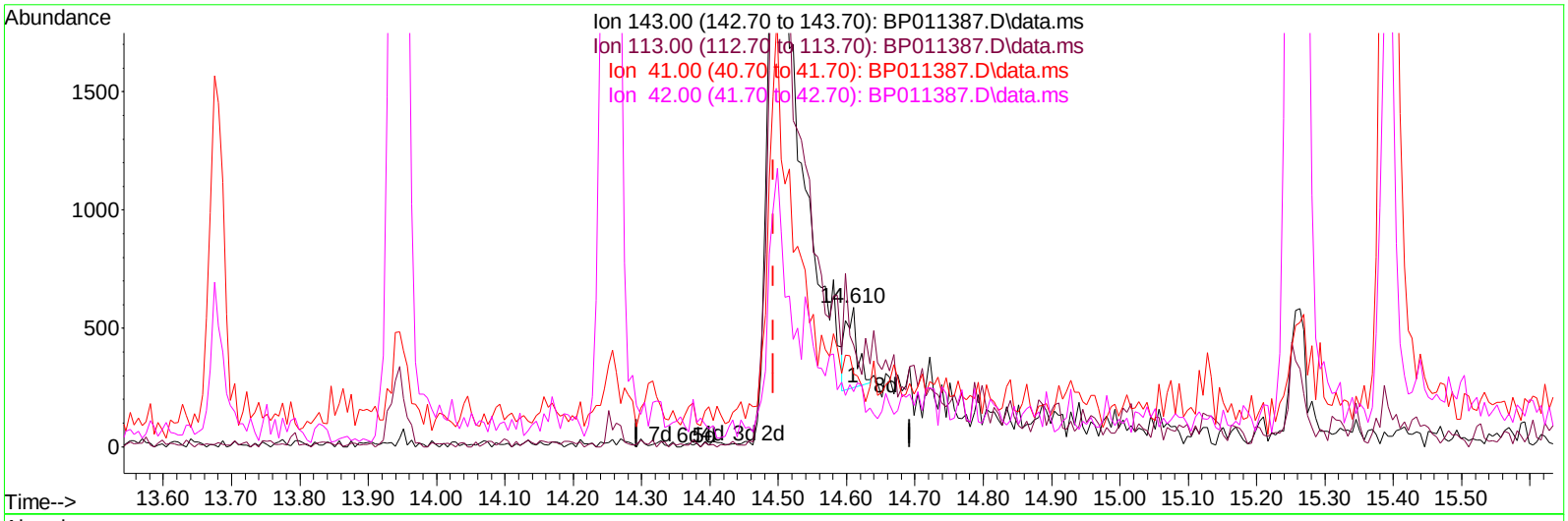

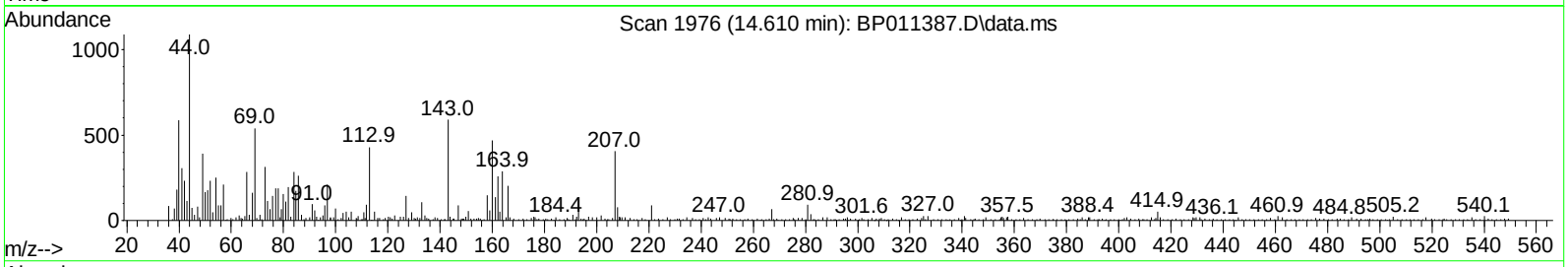

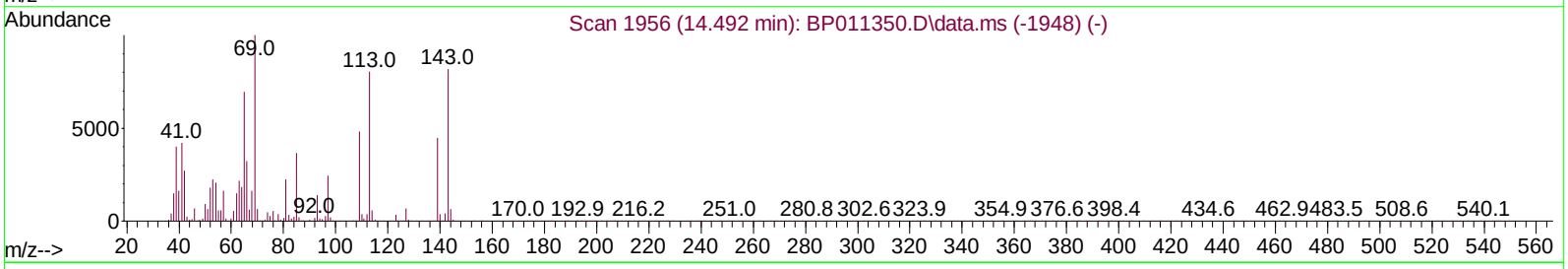

TIC: BP011387.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.610min (+ 0.118) 0.16 ng/ul

response 402

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	72.71
41.00	54.40	52.03
42.00	35.50	39.49

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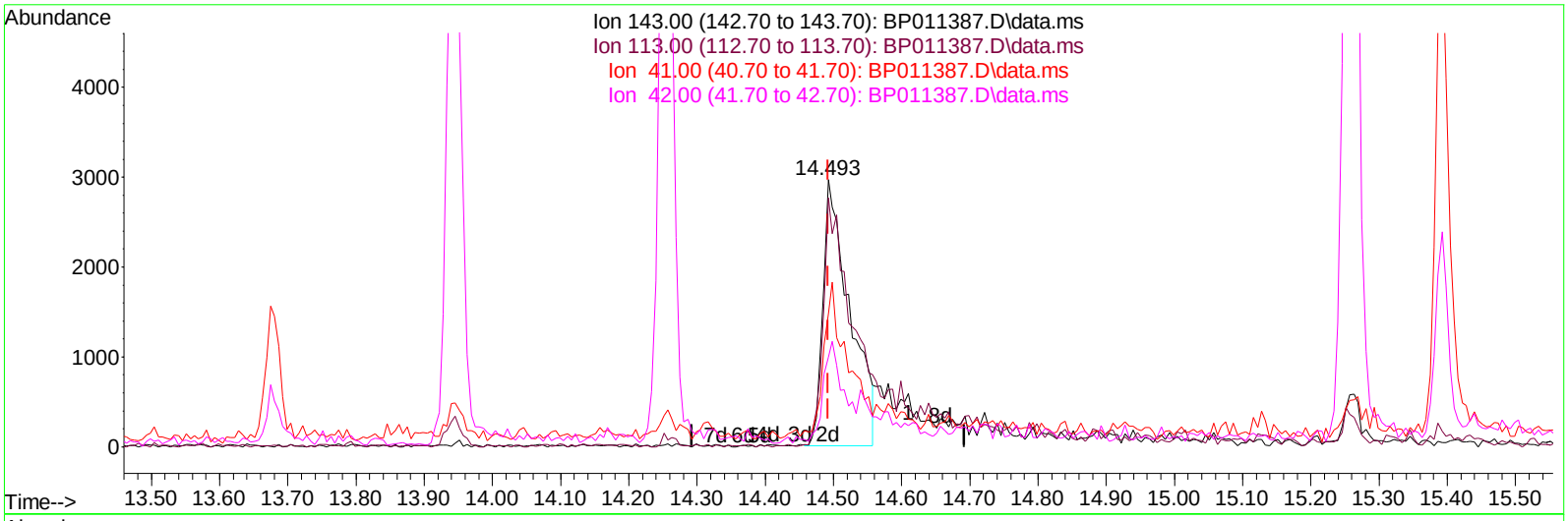

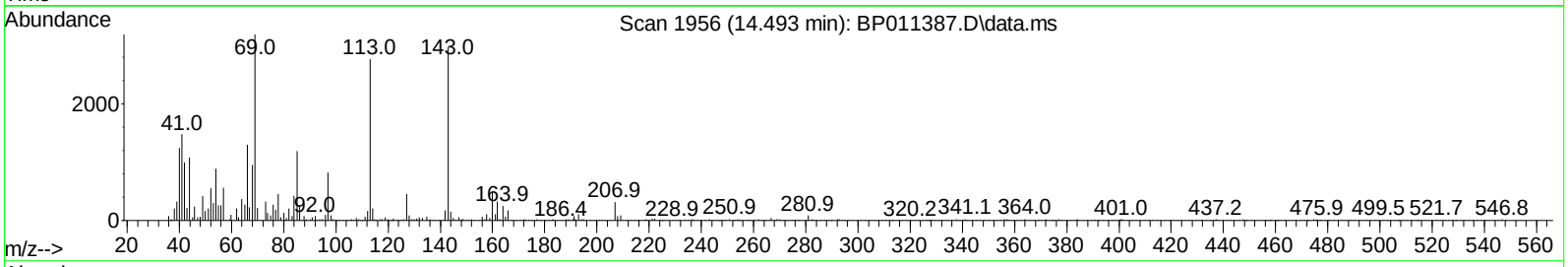

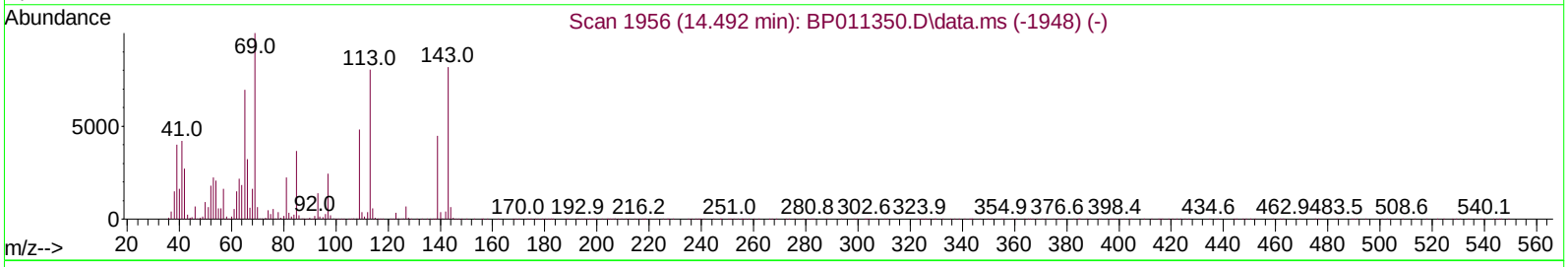

TIC: BP011387.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.000) 3.19 ng/ul m

response 8082

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.34#
41.00	54.40	49.81
42.00	35.50	33.63

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	84364	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.369	136	354950	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.257	164	199690	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	420677	20.000	ng/ul	-0.01
79) Chrysene-d12	21.163	240	408494	20.000	ng/ul	# 0.00
88) Perylene-d12	23.474	264	442459	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.081	96	11189m	4.066	ng/uL	0.00
4) Pyridine-d5	3.499	84	59103	8.514	ng/ul	0.00
7) Phenol-d5	6.764	99	46302	6.202	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	139056	26.268	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.111	132	129278	22.069	ng/ul	-0.01
15) 4-Methylphenol-d8	8.299	113	84409	14.399	ng/ul	-0.02
21) Nitrobenzene-d5	8.746	128	74572	30.870	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.463	143	71060	32.129	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.999	165	114772	25.697	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.522	131	168323	21.556	ng/ul	-0.01
46) Dimethylphthalate-d6	13.681	166	449857	32.468	ng/ul	0.00
49) Acenaphthylene-d8	13.946	160	508214	30.471	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.493	143	8082m	3.194	ng/ul	0.00
60) Fluorene-d10	15.257	176	374787	31.630	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.393	200	55956	28.637	ng/ul	0.00
73) Anthracene-d10	17.128	188	670239	35.446	ng/ul	-0.01
81) Pyrene-d10	19.392	212	789925	35.099	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.322	264	781695	35.032	ng/ul	0.00

Target Compounds Qvalue

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

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