Quantitation Report (Qedit)

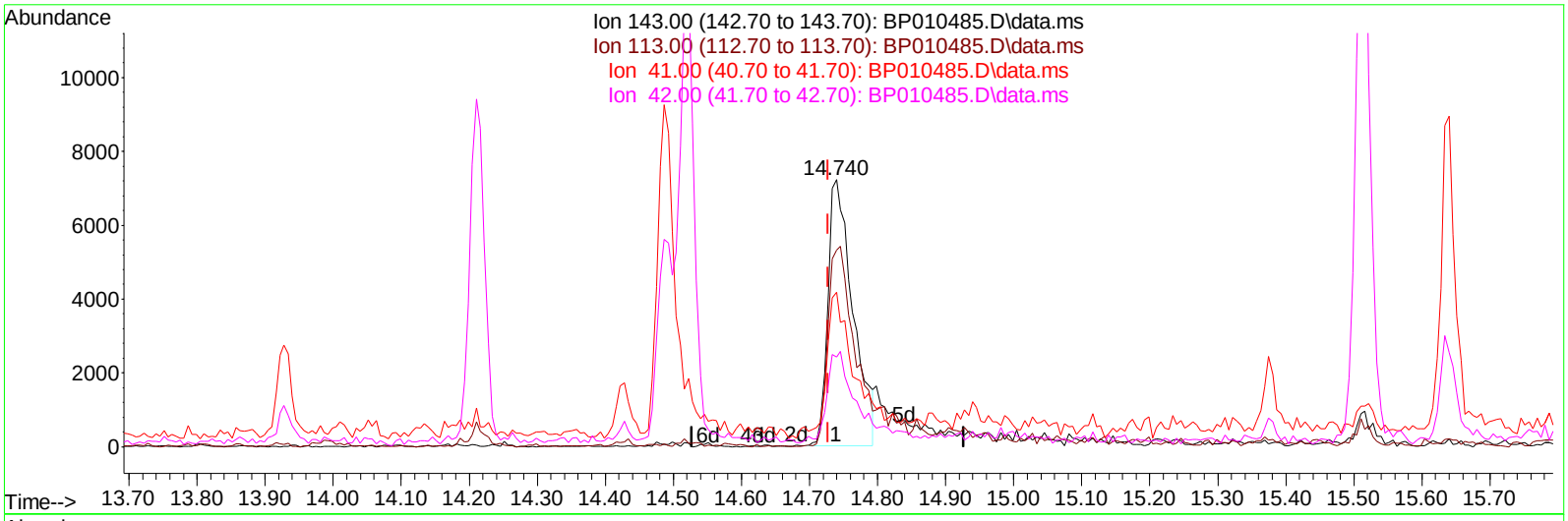
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
Data File : BP010485.D
Acq On : 23 May 2022 20:00
Operator : CG/JU
Sample : N2918-09
Misc :
ALS Vial : 14 Sample Multiplier: 1

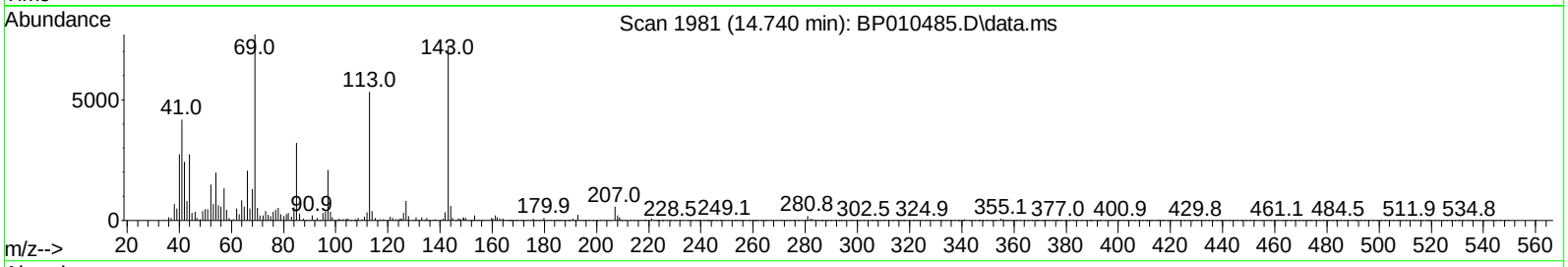
Quant Time: May 23 23:20:58 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

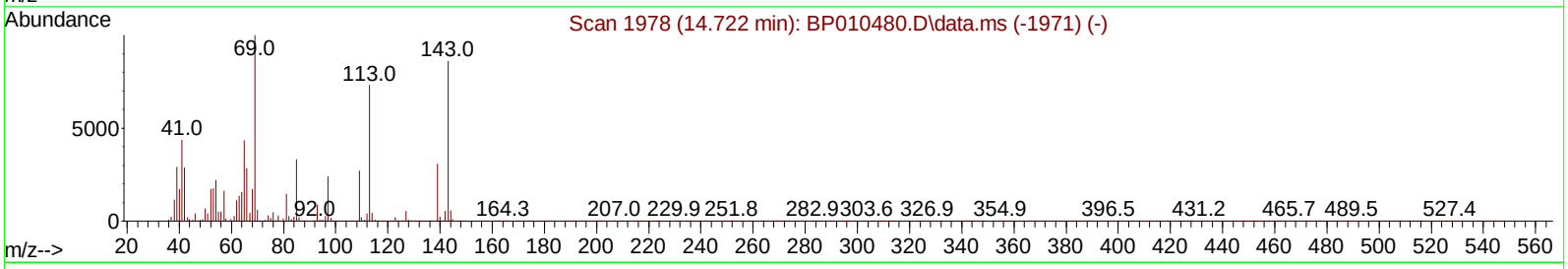
Instrument :
BNA_P
ClientSampleId :
DBTG7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/24/2022
Supervised By :Yogesh Patel 05/26/2022







TIC: BP010485.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 2.97 ng/ul

response 18331

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.72#
41.00	23.20	57.85#
42.00	18.00	33.52#

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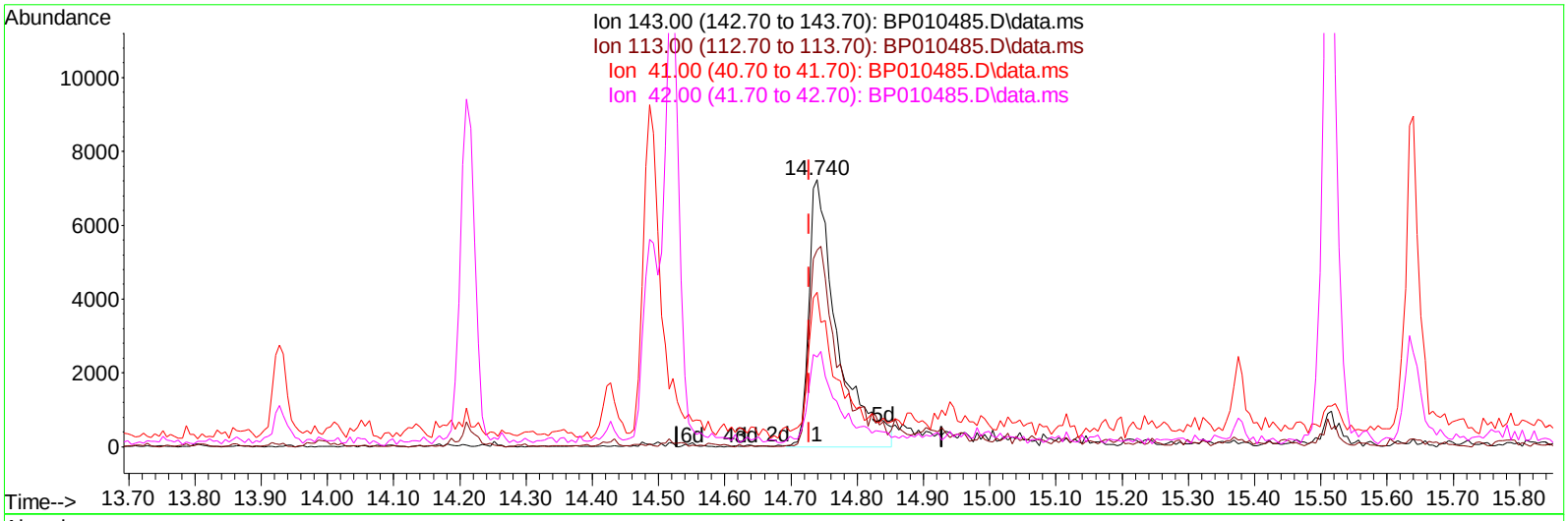

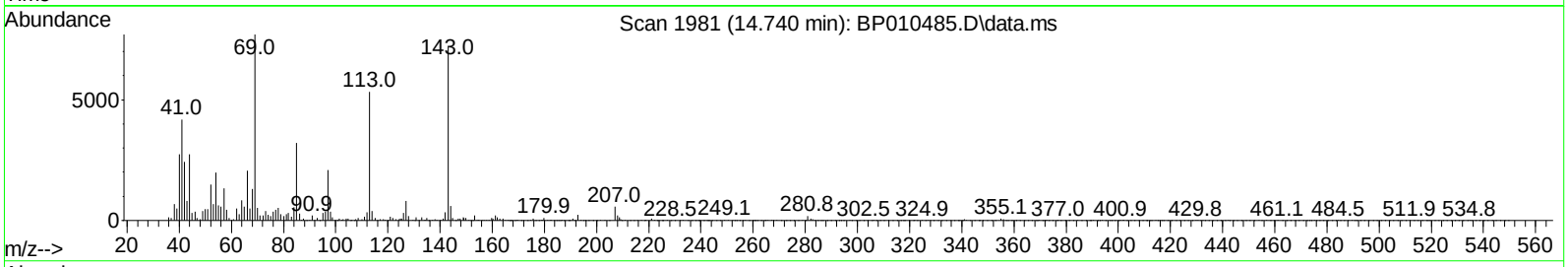

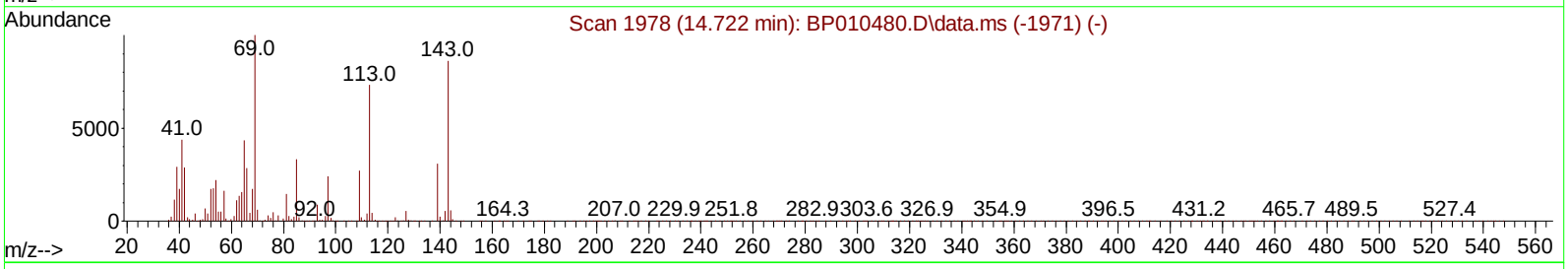

TIC: BP010485.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.50 ng/ul m

response 21622

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	73.72#
41.00	23.20	57.85#
42.00	18.00	33.52#

Quantitation Report (Qedit)

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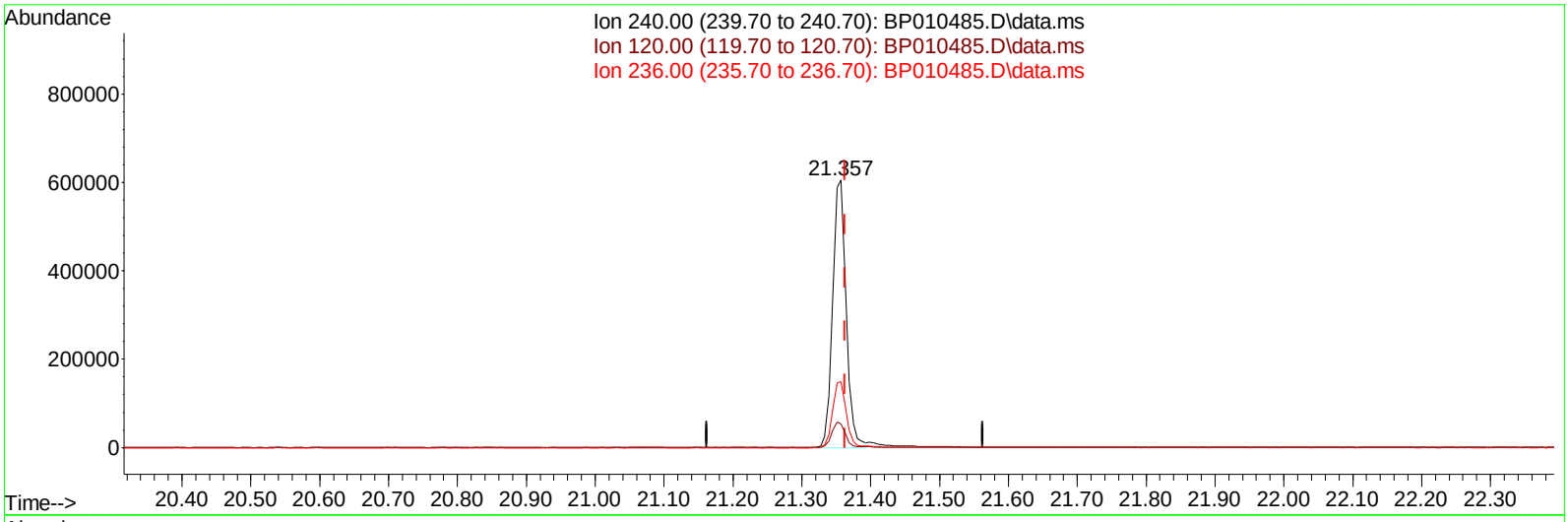

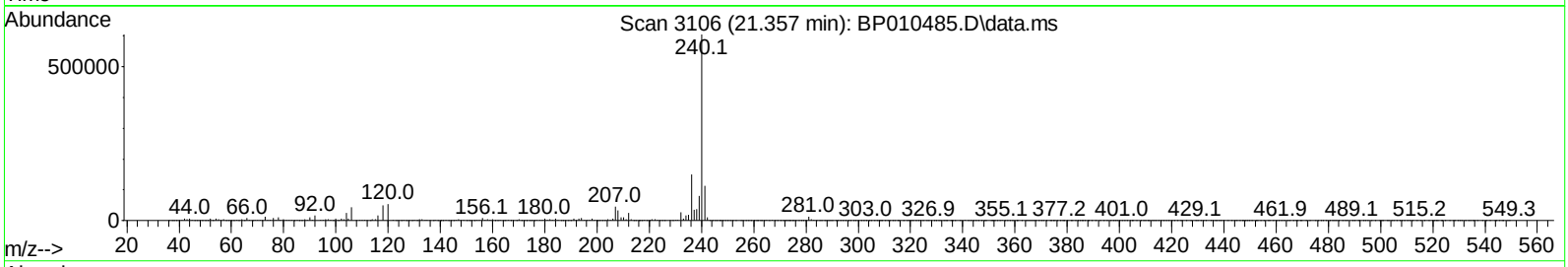

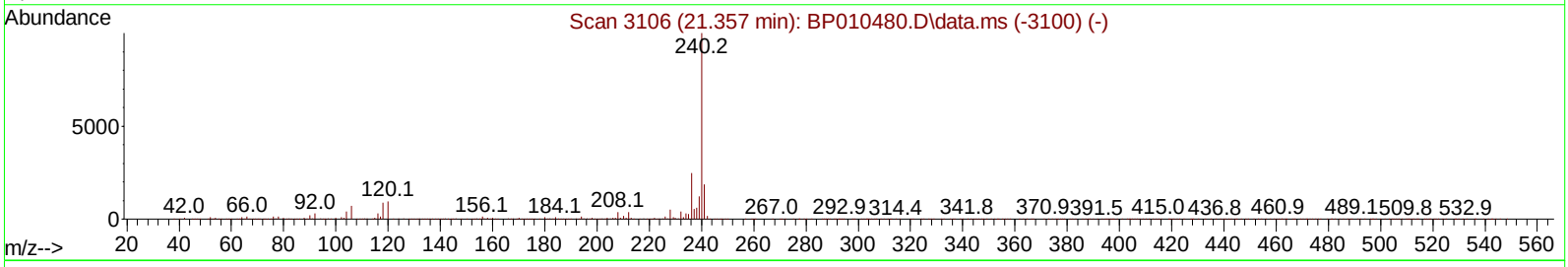

TIC: BP010485.D\data.ms

(79) Chrysene-d12 (I)

21.357min (-0.006) 20.00 ng/ul

response 810949

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.97
236.00	31.60	24.64#
0.00	0.00	0.00

Quantitation Report (Qedit)

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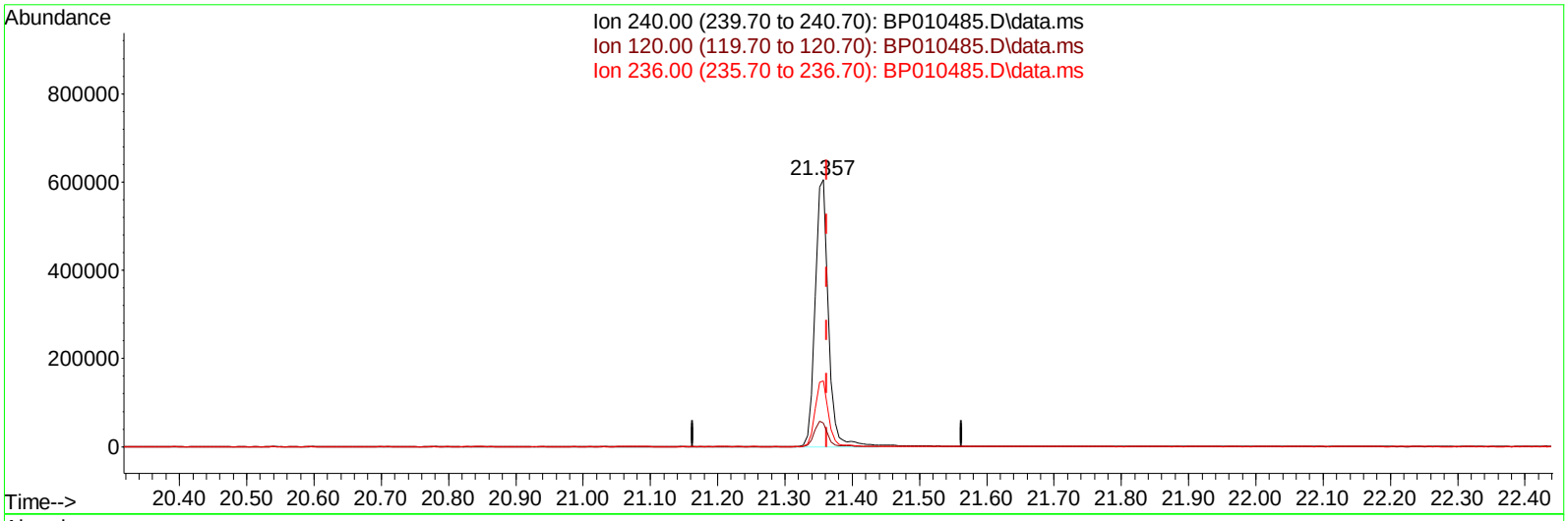

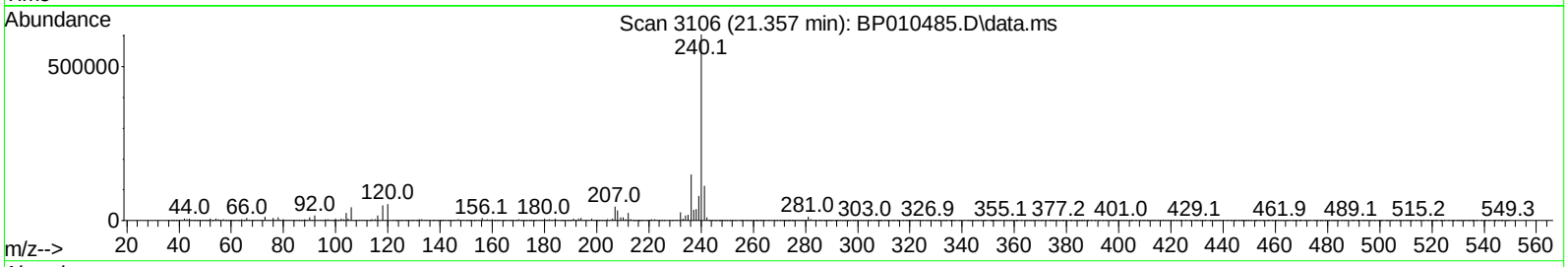

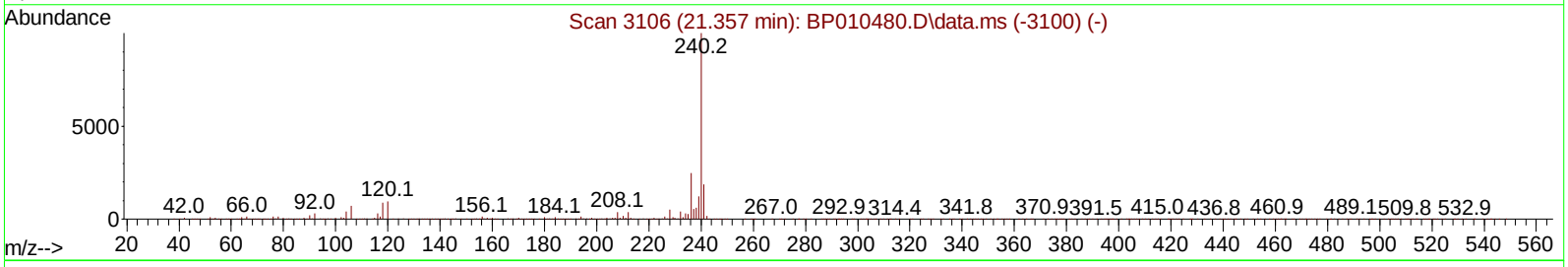

TIC: BP010485.D\data.ms

(79) Chrysene-d12 (I)

21.357min (-0.006) 20.00 ng/ul m

response 831183

Ion	Exp%	Act%
240.00	100.00	100.00
120.00	9.30	8.97
236.00	31.60	24.64#
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	7.887	152	154389	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	689042	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.522	164	471169	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	999002	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	831183m	20.000	ng/ul	0.00
88) Perylene-d12	23.774	264	661002	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	13582	3.750	ng/uL	0.00
4) Pyridine-d5	3.764	84	91304	8.878	ng/ul	0.00
7) Phenol-d5	7.058	99	64136	4.949	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	208740	23.976	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	191934	19.357	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	127132	11.544	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	130230	24.421	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	131427	23.256	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	226944	21.237	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	287459	18.092	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	885919	25.520	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	970826	24.245	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.740	143	21622m	3.499	ng/ul	0.01
60) Fluorene-d10	15.516	176	759568	25.151	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	114885	19.663	ng/ul	0.00
73) Anthracene-d10	17.375	188	1333727	29.484	ng/ul	0.00
81) Pyrene-d10	19.604	212	1588568	35.988	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.622	264	1024288	30.860	ng/ul	-0.01

Target Compounds Qvalue

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

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