Quantitation Report (Qedit)

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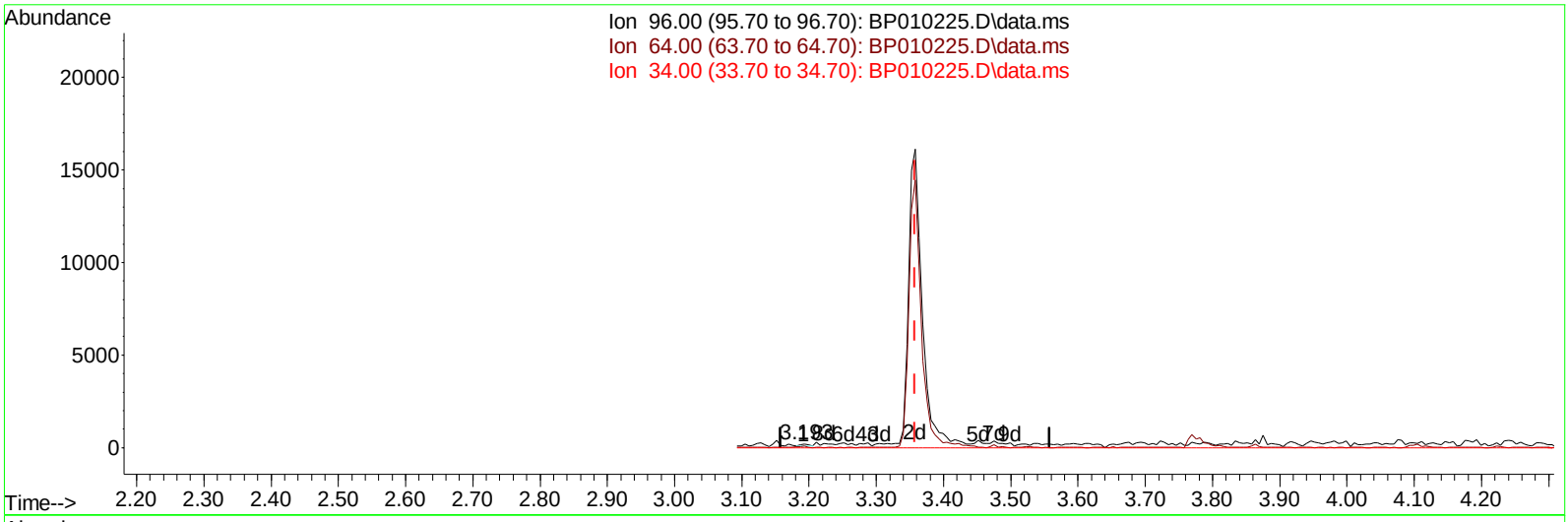
Data File : BP010225.D
Acq On : 04 May 2022 16:18
Operator : CG/JU
Sample : PB144567BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

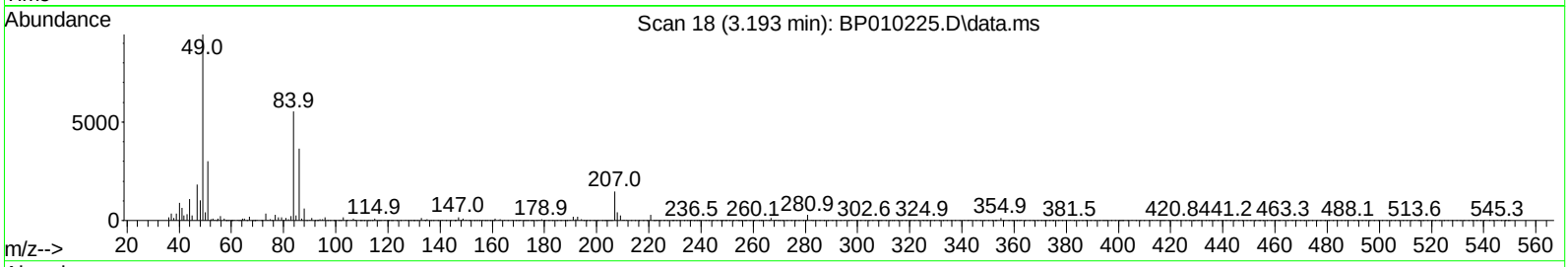
Quant Time: May 05 02:39:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

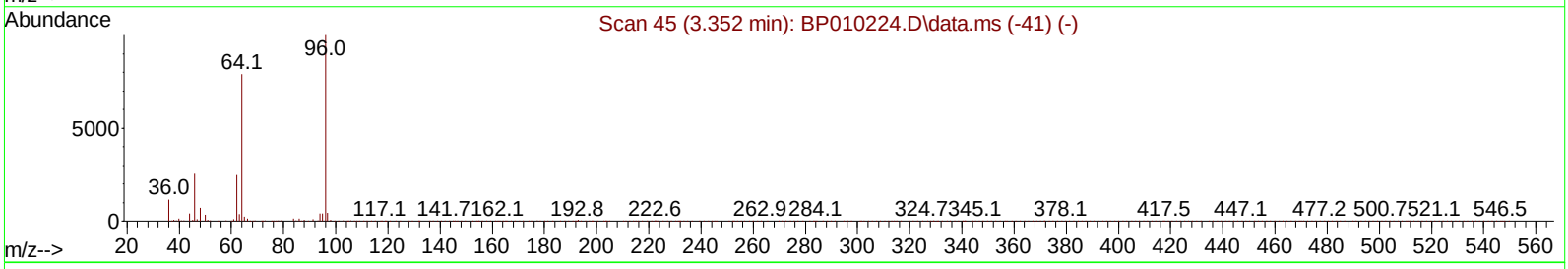
Instrument :
BNA_P
ClientSampleId :
SBLK567

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :mohammad ahmed 05/09/2022







TIC: BP010225.D\data.ms

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

3.193min (-0.165) 0.03 ng/uL

response 104

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	48.59
34.00	0.00	0.00
0.00	0.00	0.00

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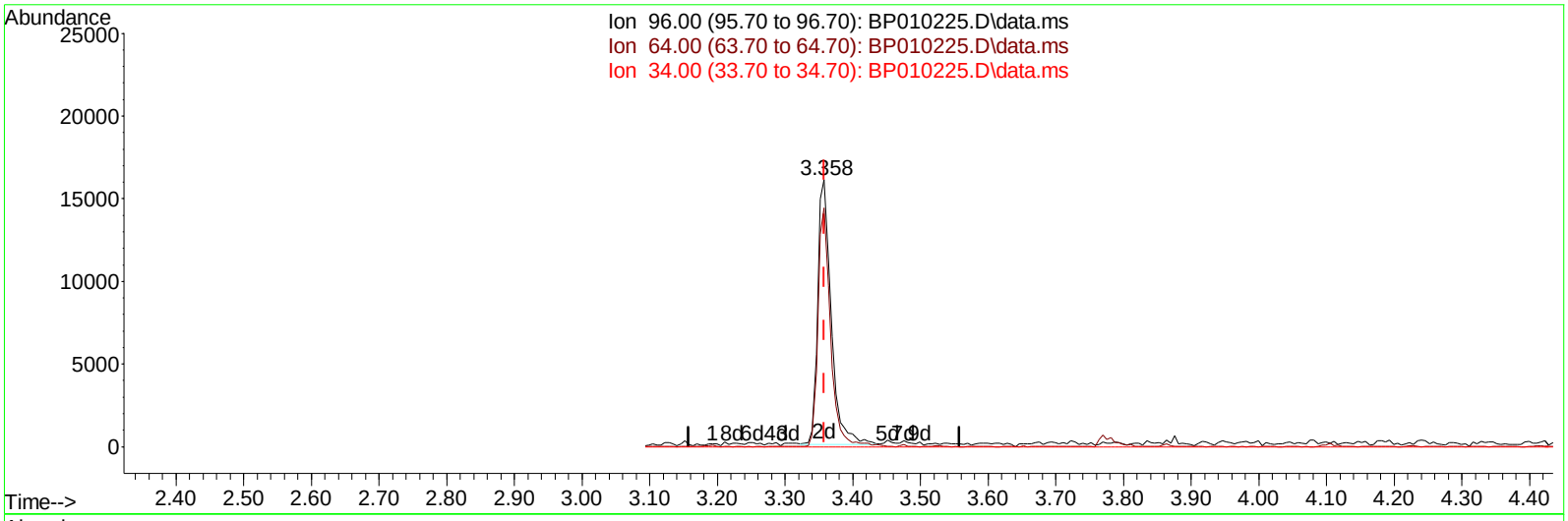

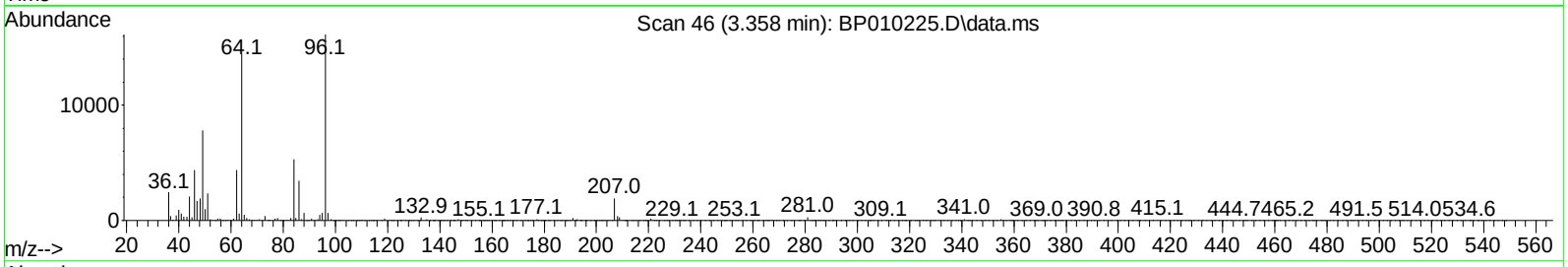

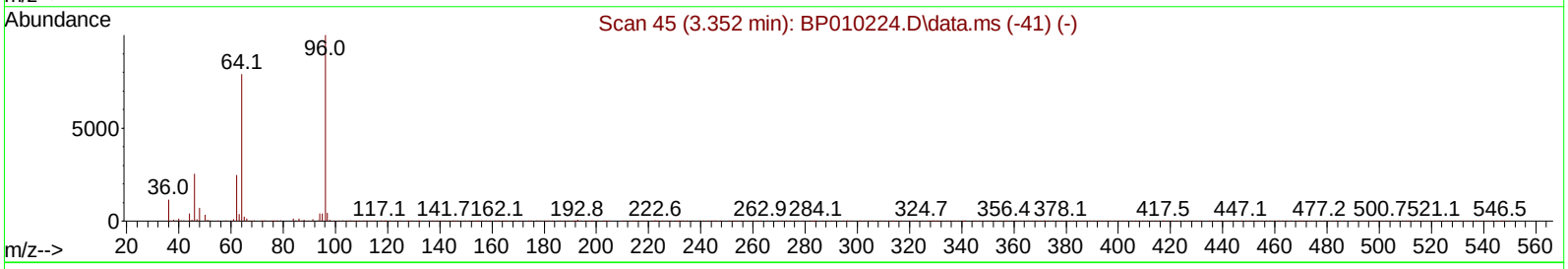

TIC: BP010225.D\data.ms

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

3.358min (-0.000) 6.08 ng/uL m

response 22130

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.50#
34.00	0.00	0.00
0.00	0.00	0.00

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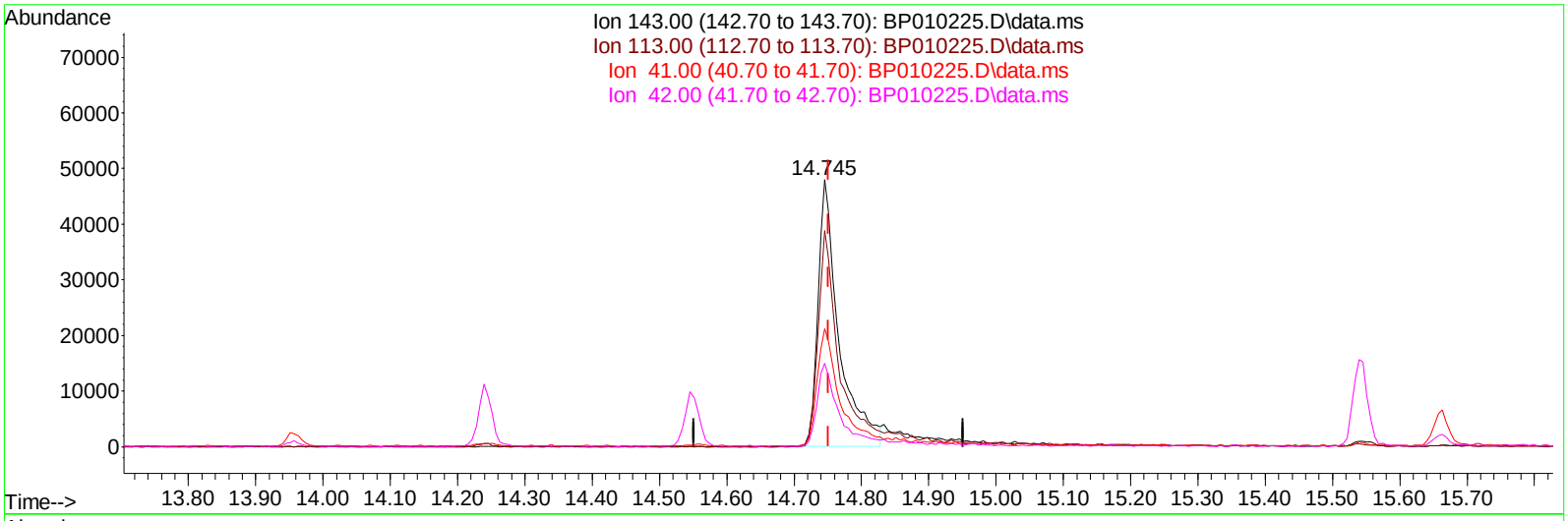

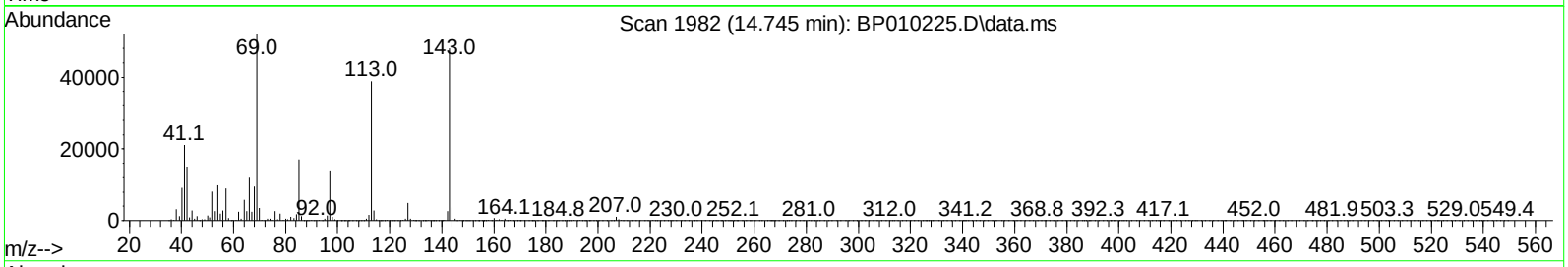

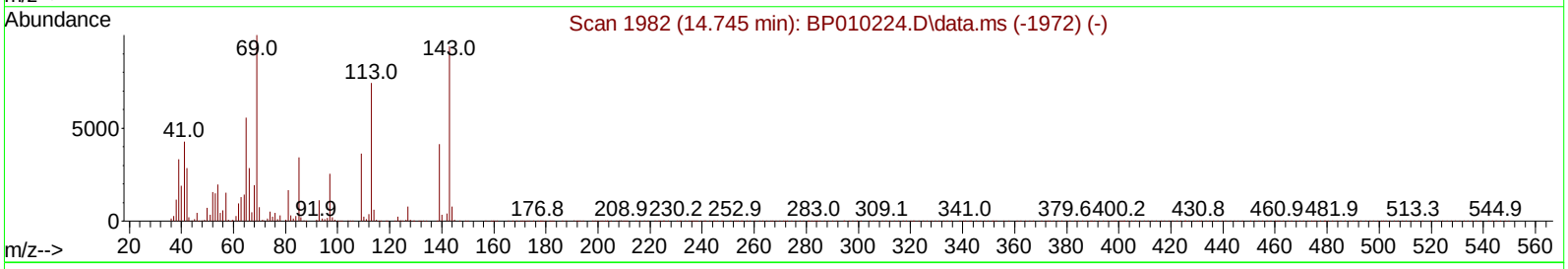

TIC: BP010225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 16.42 ng/ul

response 103616

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.07#
41.00	23.20	44.17#
42.00	18.00	31.25#

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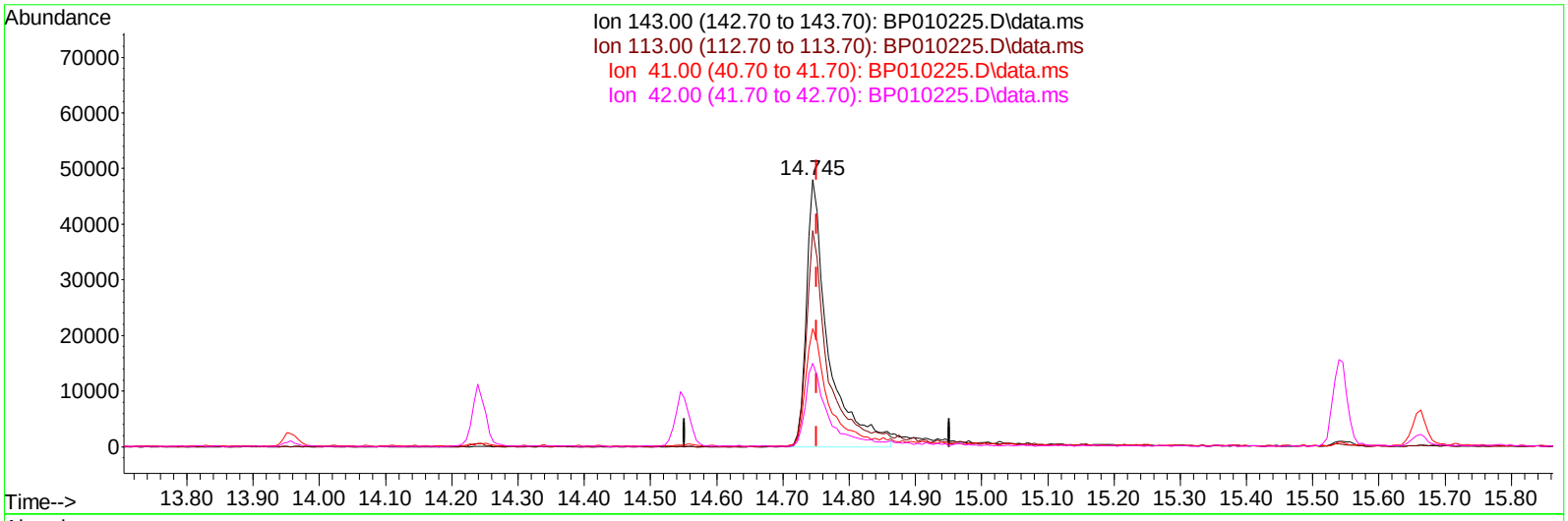

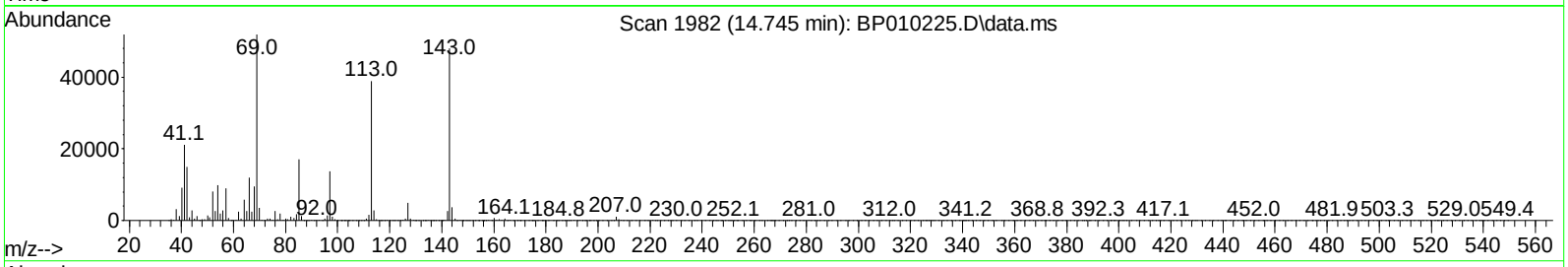

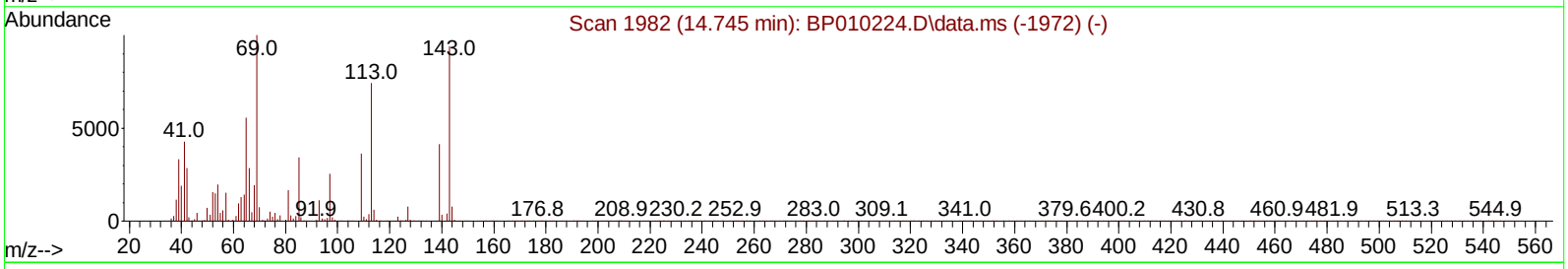

TIC: BP010225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (-0.006) 17.40 ng/ul m

response 109765

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.07#
41.00	23.20	44.17#
42.00	18.00	31.25#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	158786	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	689390	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	440615	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	785835	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	610364	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	568857	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	22130m	6.077	ng/uL	0.00
4) Pyridine-d5	3.769	84	311984	30.456	ng/ul	0.00
7) Phenol-d5	7.081	99	392907	27.204	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	280617	32.380	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	321828	30.088	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	333199	27.844	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	171002	31.187	ng/ul	0.00
24) 2-Nitrophenol-d4	9.799	143	183742	30.684	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	310327	26.620	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	456646	27.319	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1057223	30.845	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1196387	28.841	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.745	143	109765m	17.396	ng/ul	0.00
60) Fluorene-d10	15.545	176	867676	29.429	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	103235	20.776	ng/ul	0.00
73) Anthracene-d10	17.404	188	1218155	32.042	ng/ul	0.00
81) Pyrene-d10	19.633	212	1244566	35.854	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1021819	34.263	ng/ul	-0.01

Target Compounds Qvalue

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

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