Quantitation Report (Qedit)

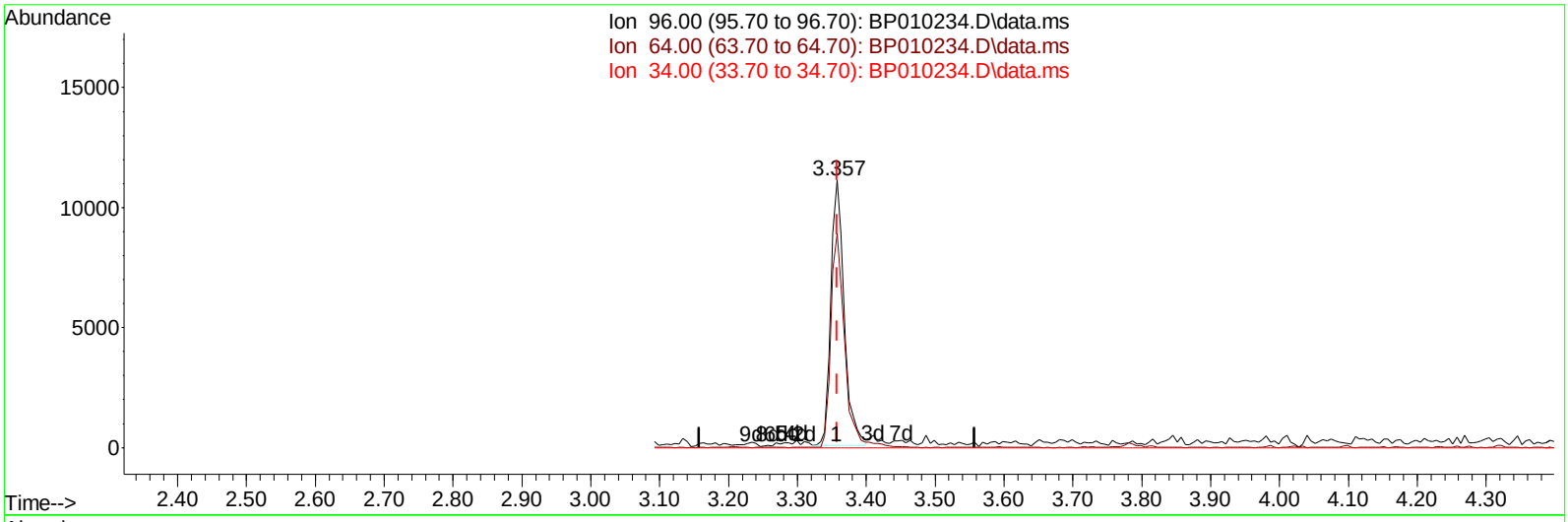
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050422\
Data File : BP010234.D
Acq On : 04 May 2022 21:22
Operator : CG/JU
Sample : N2668-10
Misc :
ALS Vial : 12 Sample Multiplier: 1

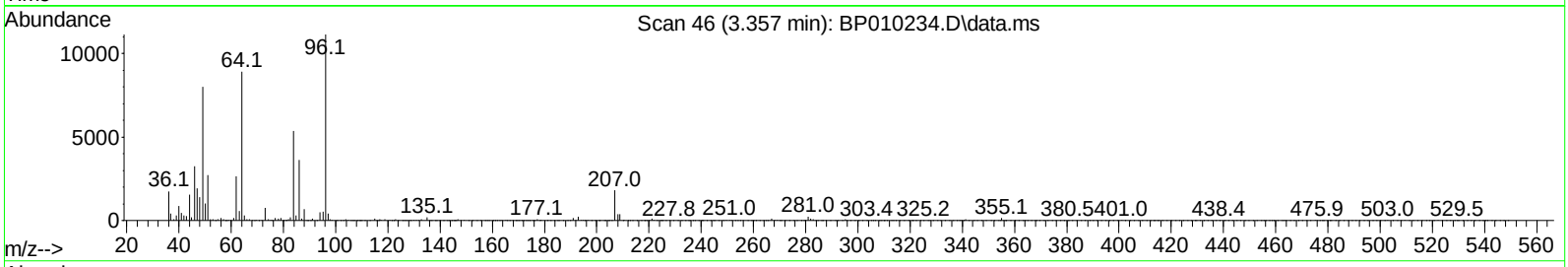
Quant Time: May 05 02:41:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

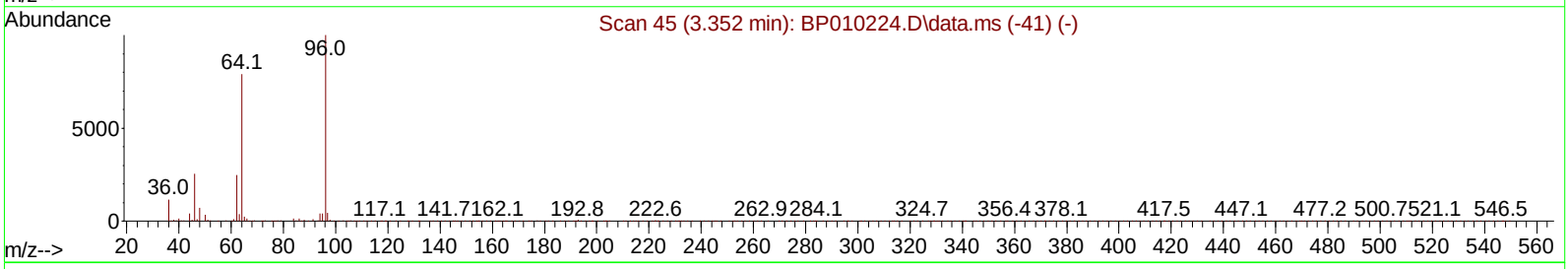
Instrument :
BNA_P
ClientSampleId :
EXL75

Manual IntegrationsAPPROVED

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







TIC: BP010234.D\data.ms

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

3.357min (-0.000) 4.18 ng/uL

response 14783

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.13#
34.00	0.00	0.00
0.00	0.00	0.00

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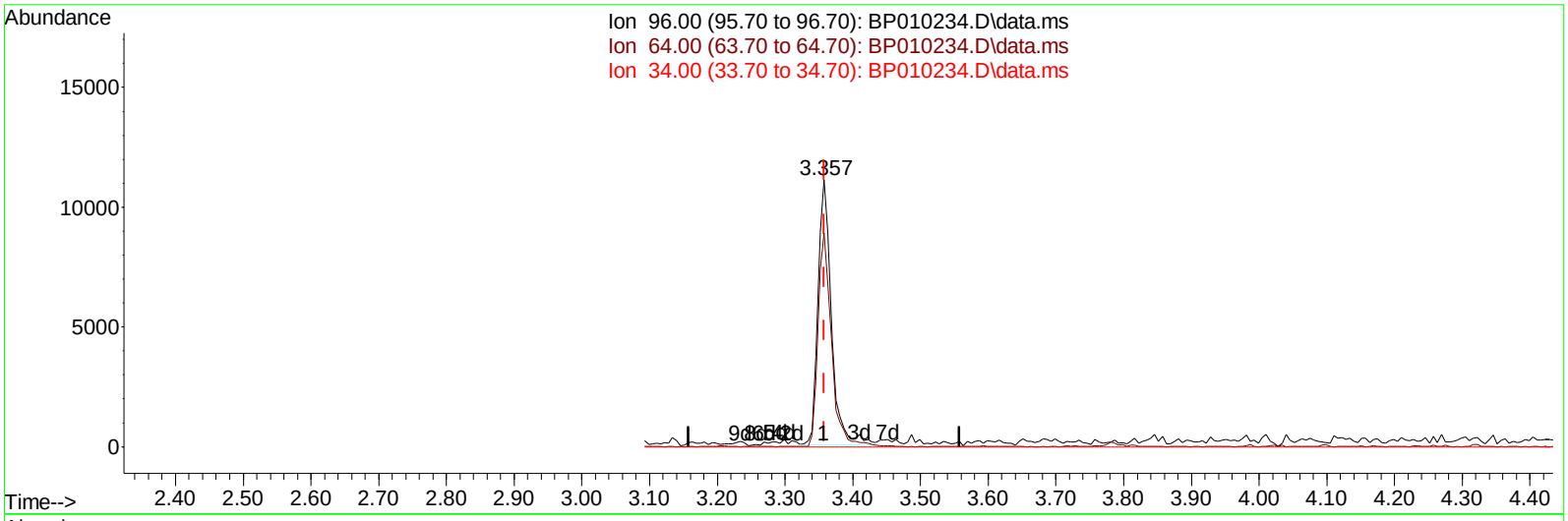

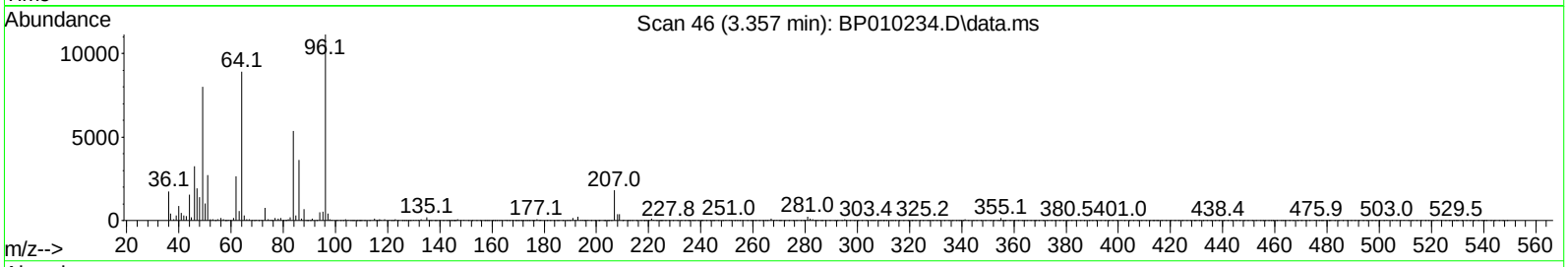

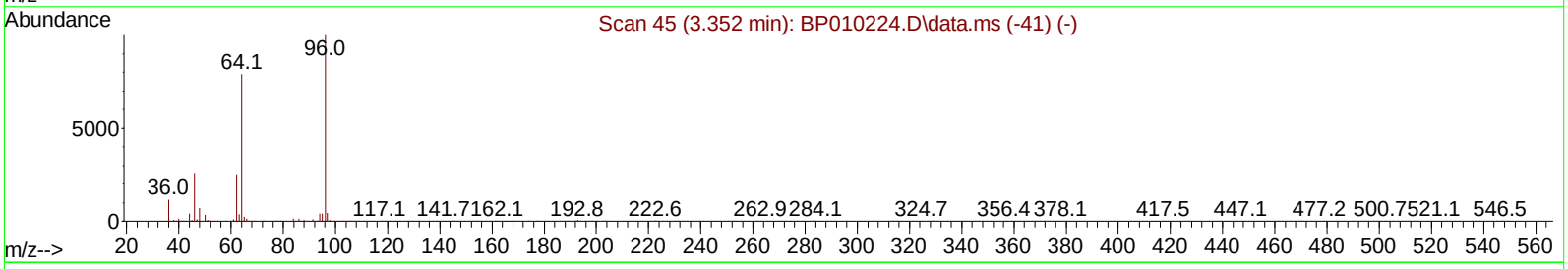

TIC: BP010234.D\data.ms

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

3.357min (-0.000) 4.33 ng/uL m

response 15330

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	80.13#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
EXL75

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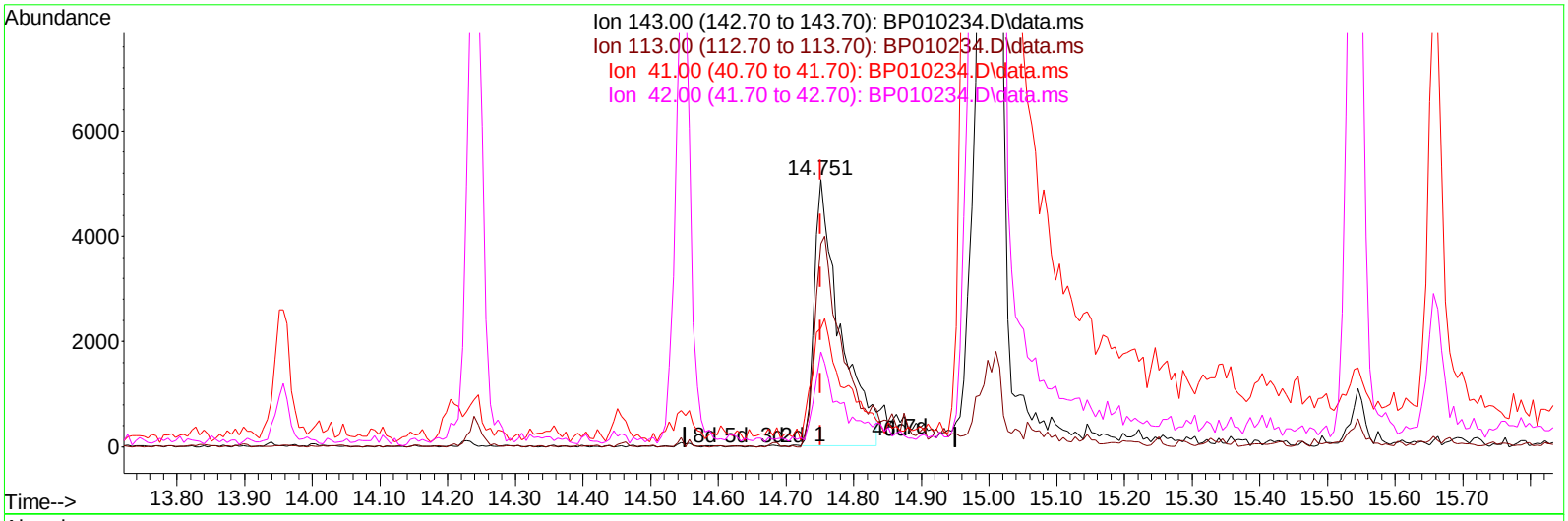

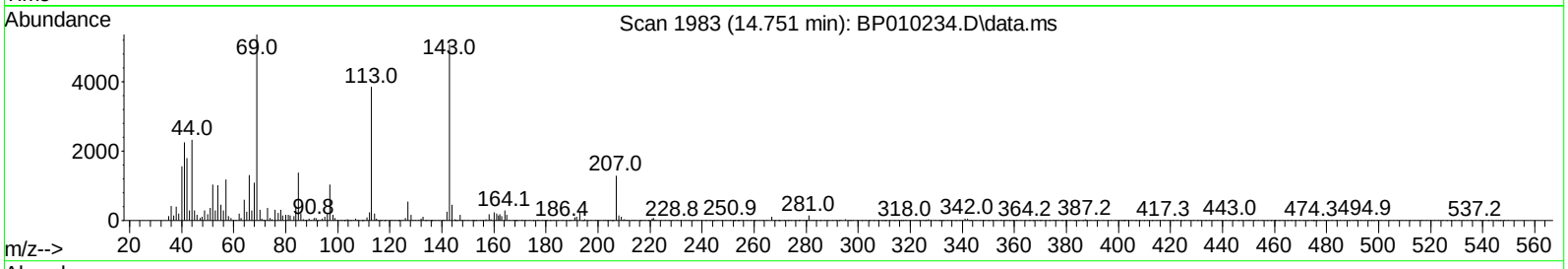

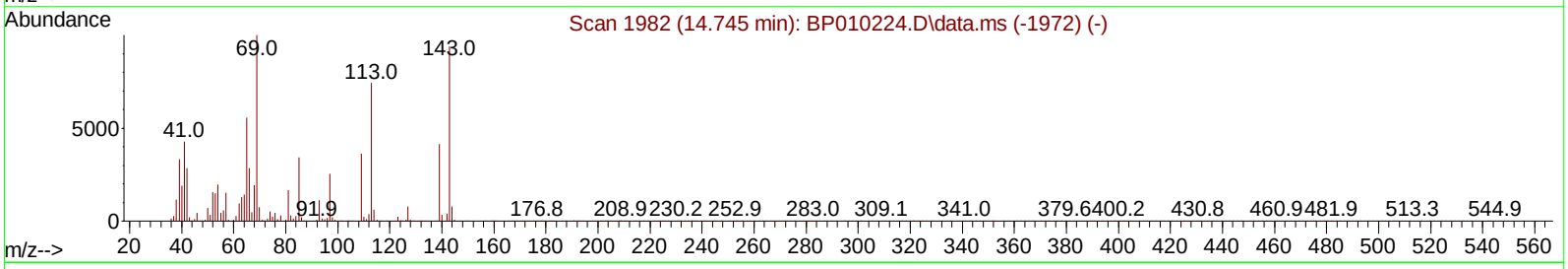

TIC: BP010234.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 2.23 ng/ul

response 13522

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.99#
41.00	23.20	44.39#
42.00	18.00	35.58#

Quantitation Report (Qedit)

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ClientSampleId :
EXL75

Manual IntegrationsAPPROVED

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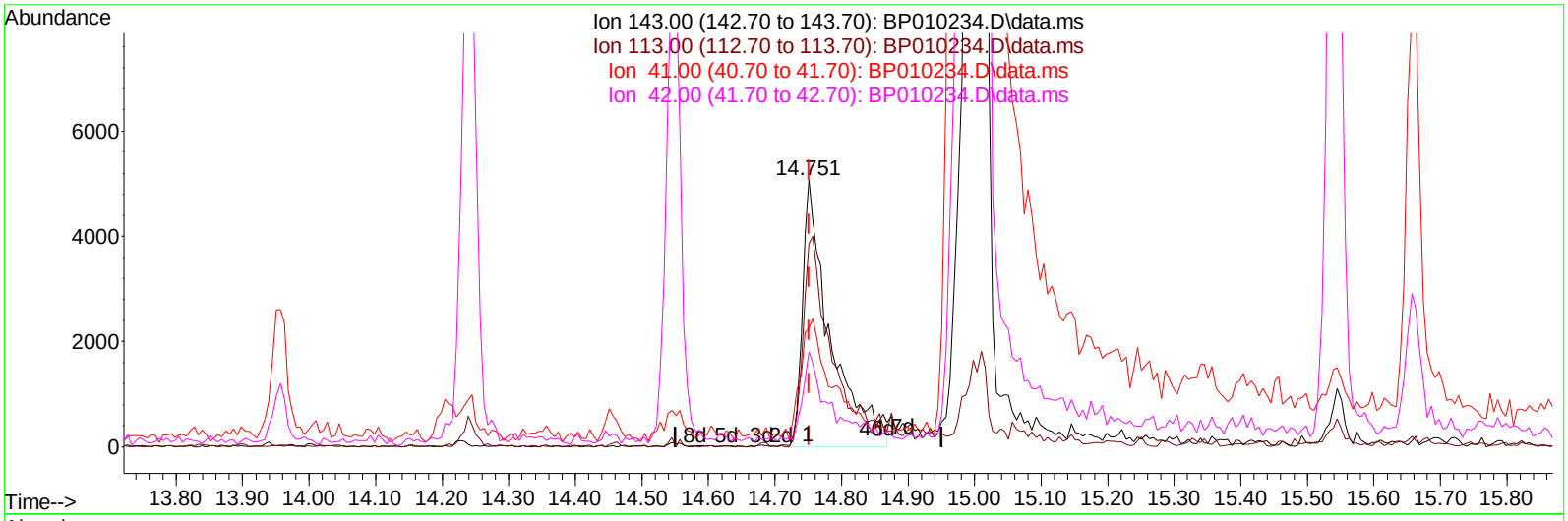
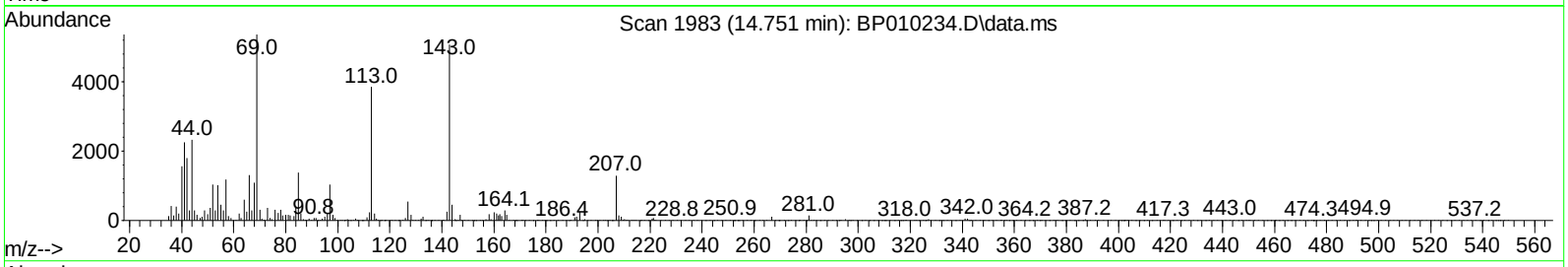
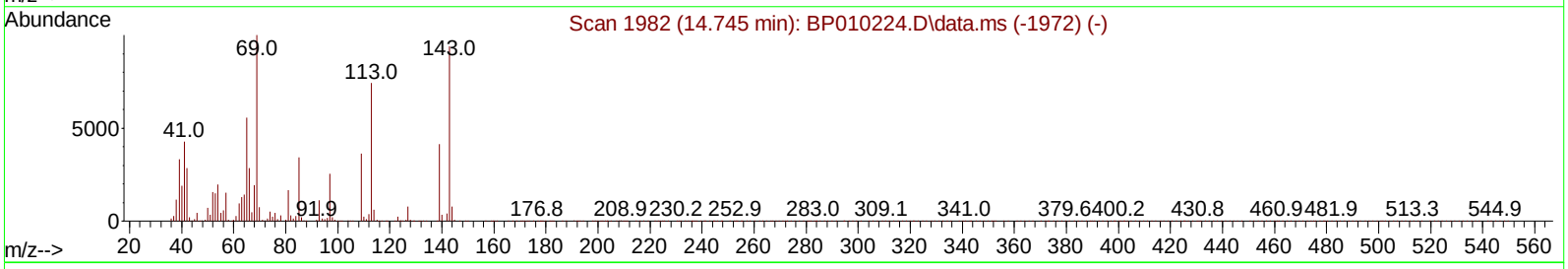
Supervised By :mohammad ahmed 05/09/2022







TIC: BP010234.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 2.45 ng/ul m

response 14839

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.99#
41.00	23.20	44.39#
42.00	18.00	35.58#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	154393	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	663671	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	422561	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.298	188	761260	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.380	240	581870	20.000	ng/ul	#-0.01
88) Perylene-d12	23.827	264	571477	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	15330m	4.330	ng/uL	0.00
4) Pyridine-d5	3.781	84	62857	6.311	ng/ul	0.00
7) Phenol-d5	7.081	99	64719	4.608	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.245	67	253170	30.044	ng/ul	0.00
11) 2-Chlorophenol-d4	7.445	132	219699	21.124	ng/ul	-0.01
15) 4-Methylphenol-d8	8.622	113	143058	12.295	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	152204	28.835	ng/ul	0.00
24) 2-Nitrophenol-d4	9.798	143	156928	27.222	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	276364	24.626	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	348602	21.664	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1053258	32.043	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1139474	28.643	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143	14839m	2.452	ng/ul	0.00
60) Fluorene-d10	15.539	176	860653	30.438	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.663	200	115552	24.006	ng/ul	0.00
73) Anthracene-d10	17.398	188	1220707	33.145	ng/ul	-0.01
81) Pyrene-d10	19.633	212	1246542	37.670	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1042375	34.792	ng/ul	-0.01
Target Compounds						Qvalue

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

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