Quantitation Report (Qedit)

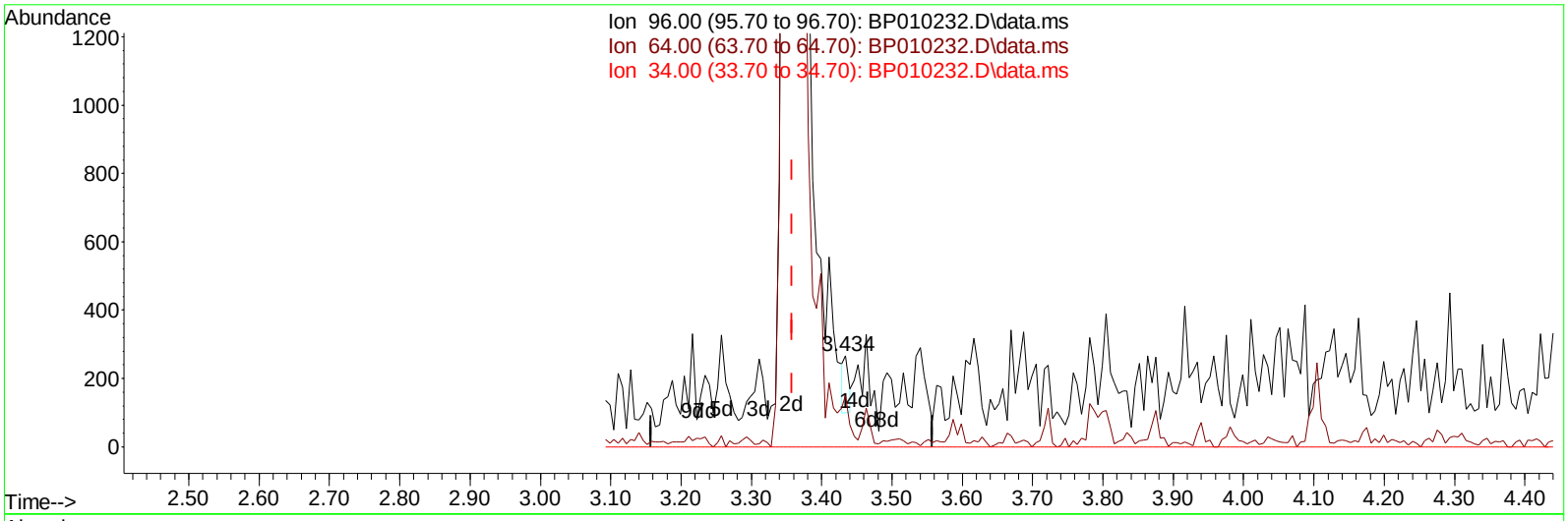
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050422\
Data File : BP010232.D
Acq On : 04 May 2022 20:15
Operator : CG/JU
Sample : N2668-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

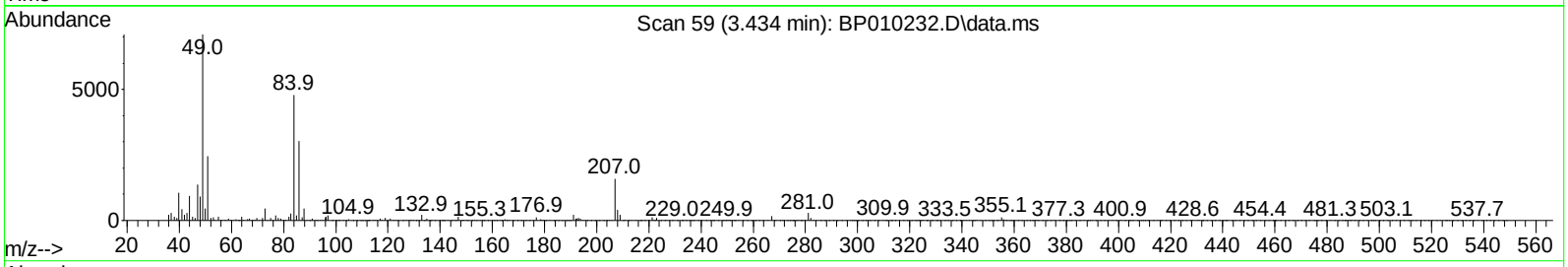
Quant Time: May 05 02:41:22 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 00:58:32 2022
Response via : Initial Calibration

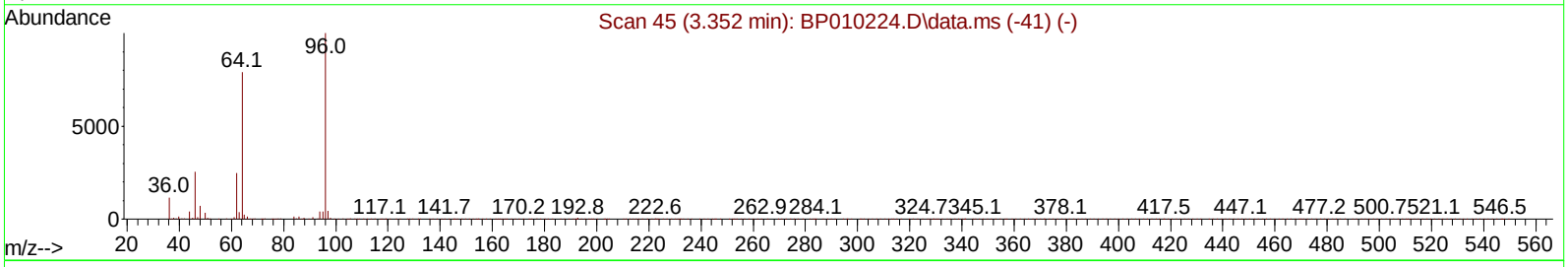
Instrument :
BNA_P
ClientSampleId :
EXL46

Manual IntegrationsAPPROVED

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







TIC: BP010232.D\data.ms

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

3.434min (+ 0.076) 0.02 ng/uL

response 83

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	64.04
34.00	0.00	0.00
0.00	0.00	0.00

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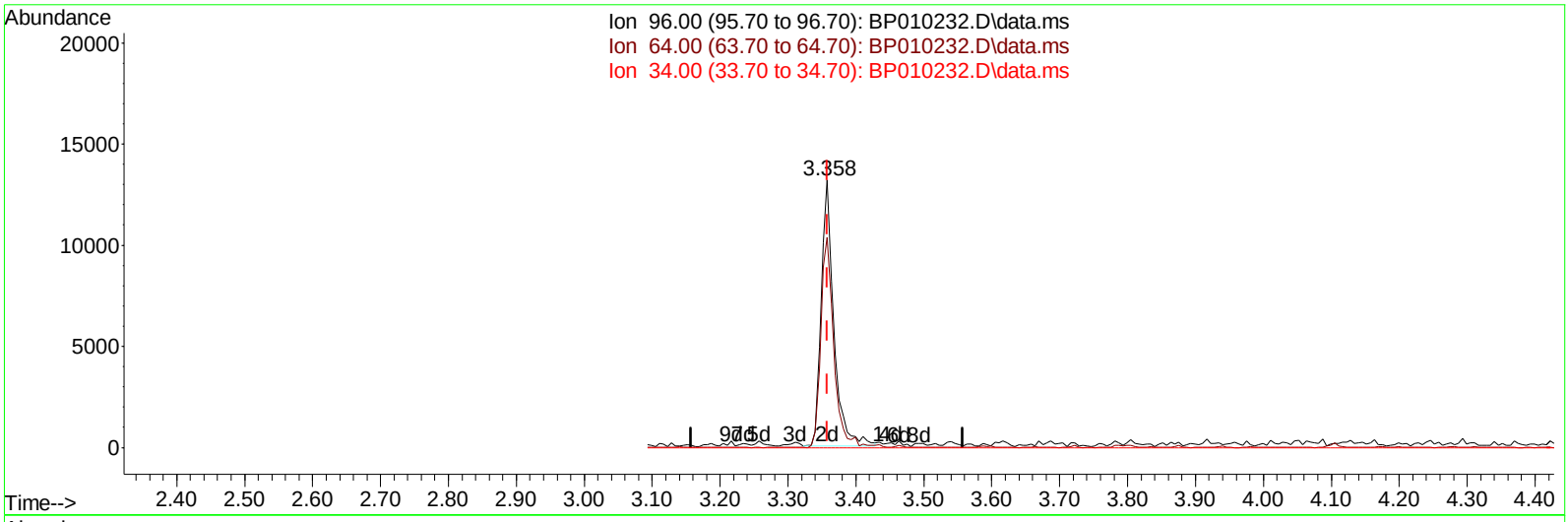

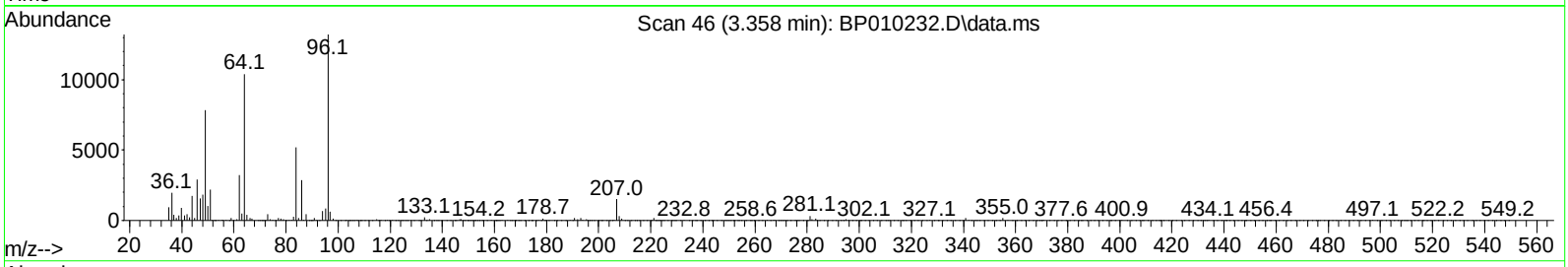

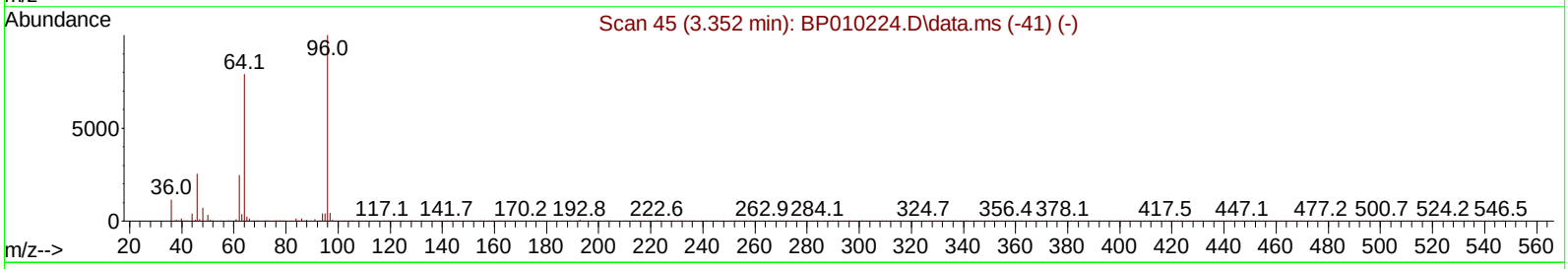

TIC: BP010232.D\data.ms

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

3.358min (-0.000) 4.30 ng/uL m

response 17064

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	78.65#
34.00	0.00	0.00
0.00	0.00	0.00

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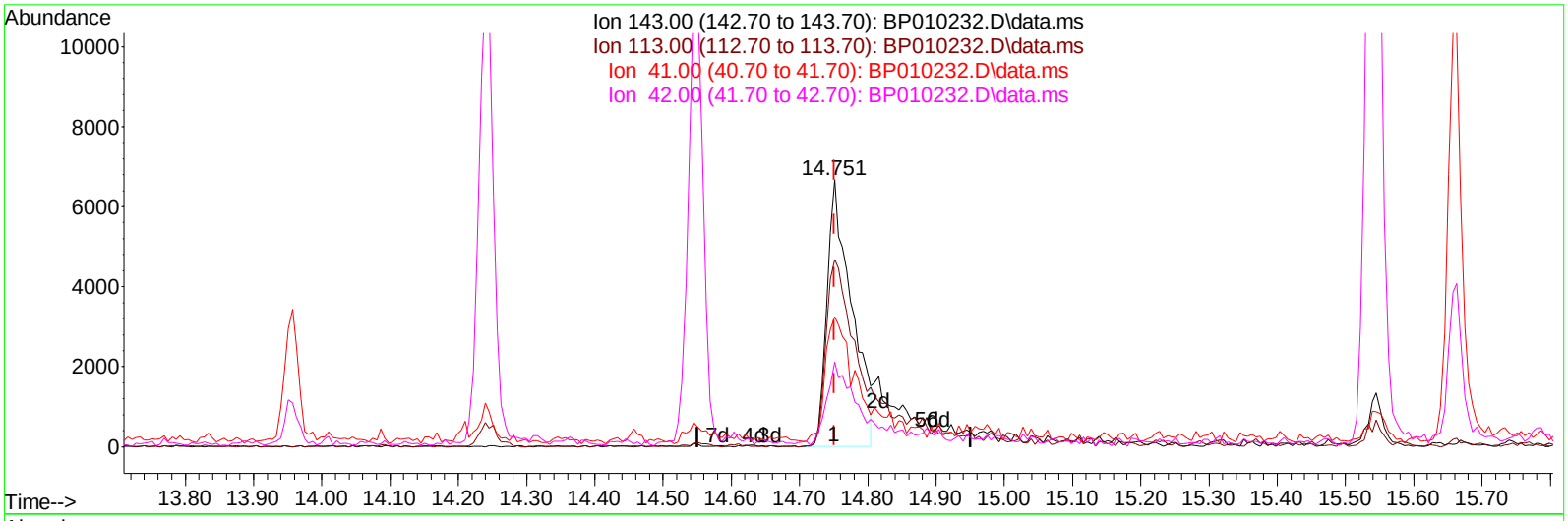

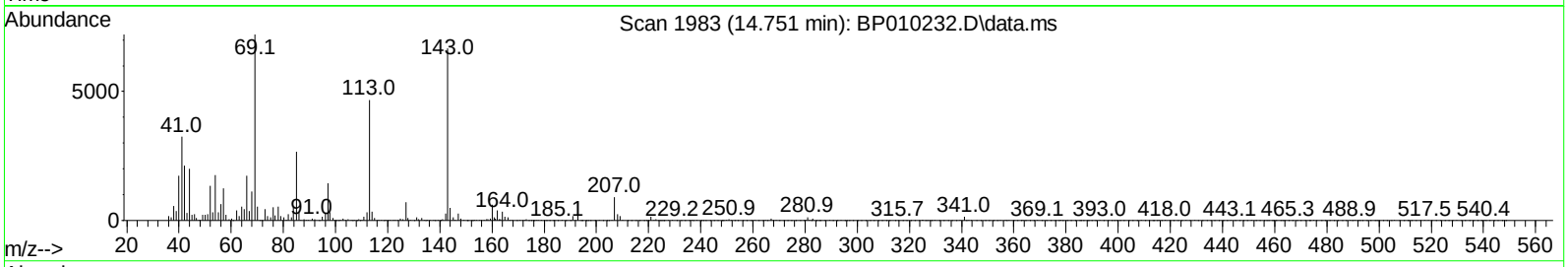

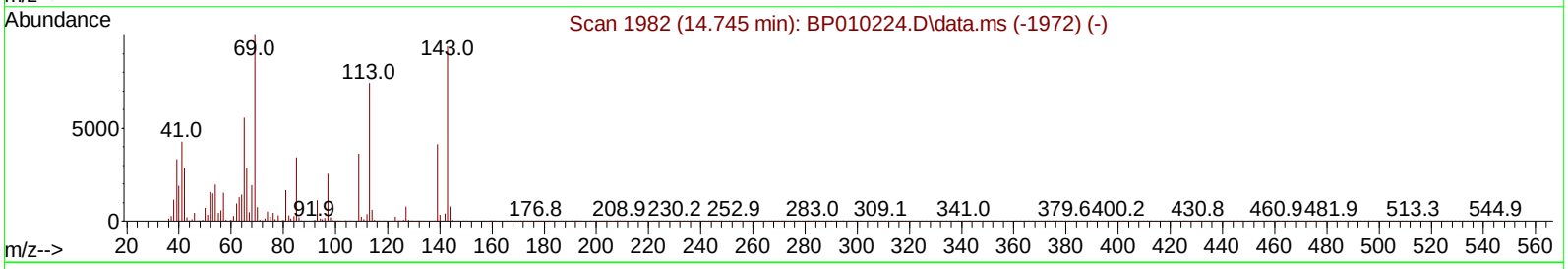

TIC: BP010232.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 2.28 ng/ul

response 16641

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	70.13#
41.00	23.20	48.82#
42.00	18.00	31.83#

Quantitation Report (Qedit)

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Manual IntegrationsAPPROVED

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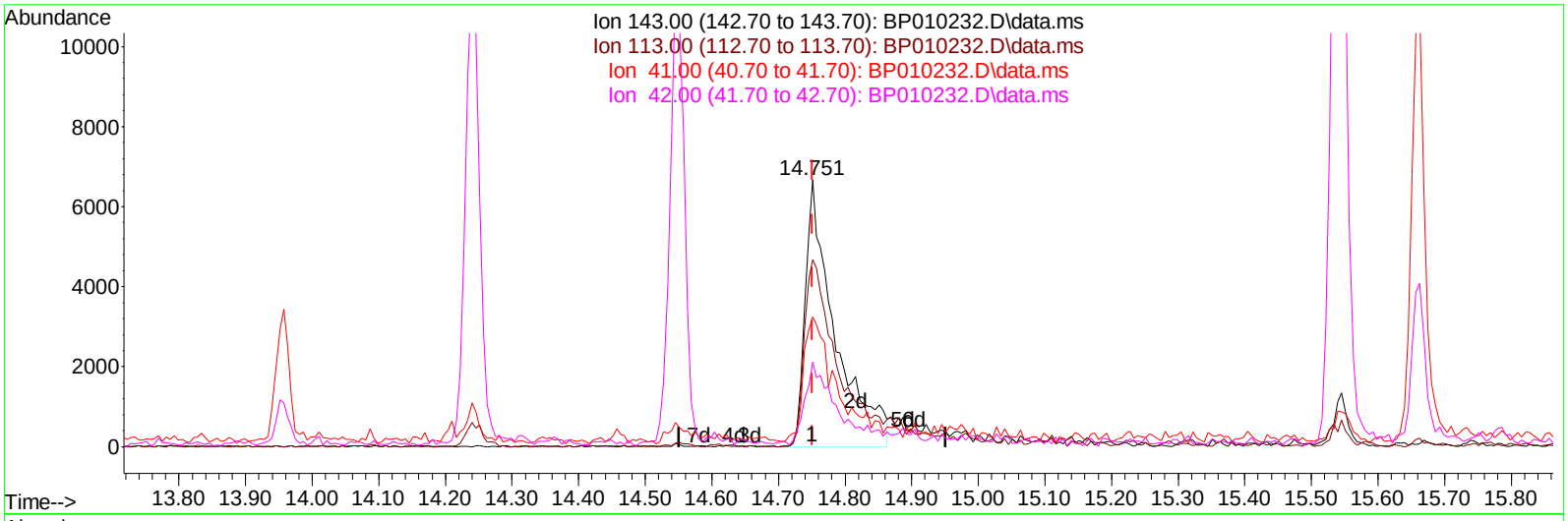

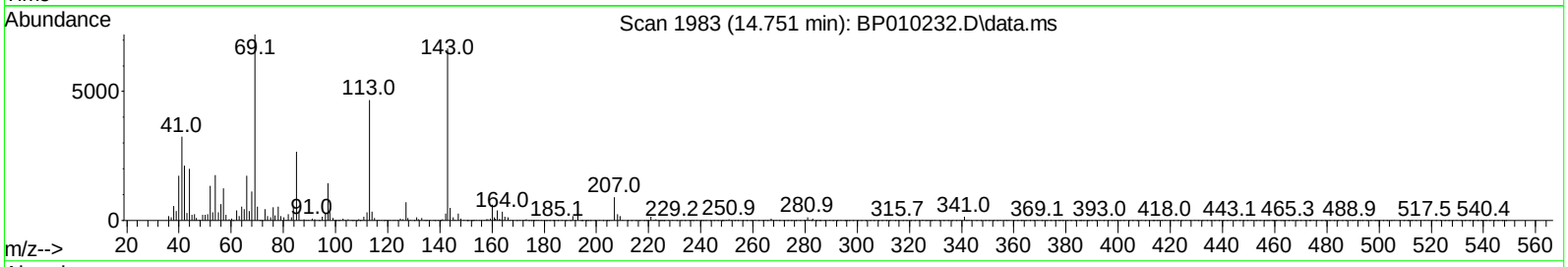

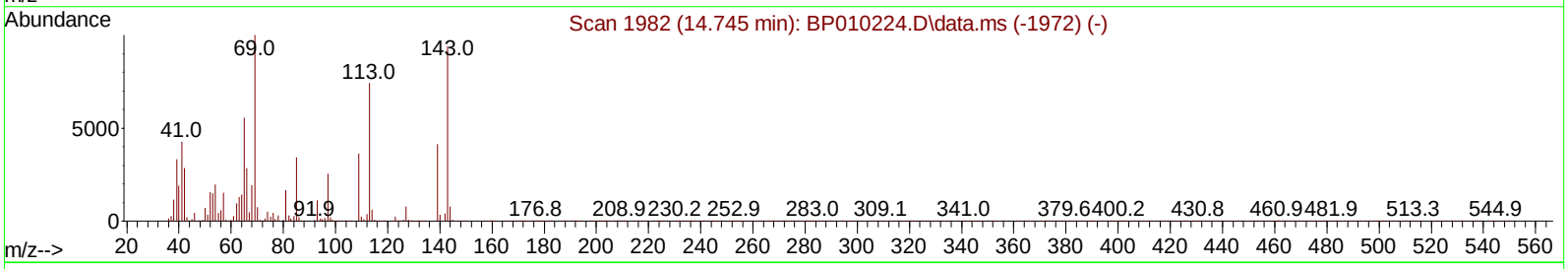

TIC: BP010232.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.000) 2.82 ng/ul m

response 20618

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	70.13#
41.00	23.20	48.82#
42.00	18.00	31.83#

Quantitation Report (QT Reviewed)

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EXL46

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	172990	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	756061	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	509881	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188	1012840	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.386	240	699081	20.000	ng/ul	# 0.00
88) Perylene-d12	23.827	264	638437	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	17064m	4.301	ng/uL	0.00
4) Pyridine-d5	3.787	84	43601	3.907	ng/ul	0.01
7) Phenol-d5	7.081	99	67014	4.259	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	279823	29.637	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	235079	20.173	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	144691	11.099	ng/ul	-0.01
21) Nitrobenzene-d5	9.075	128	174979	29.099	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	175568	26.734	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.340	165	280749	21.959	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.857	131	386711	21.095	ng/ul	0.00
46) Dimethylphthalate-d6	13.957	166	1250060	31.517	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1336203	27.836	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143	20618m	2.824	ng/ul	0.00
60) Fluorene-d10	15.545	176	1056536	30.967	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200	160530	25.066	ng/ul	0.00
73) Anthracene-d10	17.404	188	1662858	33.936	ng/ul	0.00
81) Pyrene-d10	19.633	212	1724017	43.364	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1221076	36.482	ng/ul	-0.01
Target Compounds						Qvalue

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

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