Quantitation Report (Qedit)

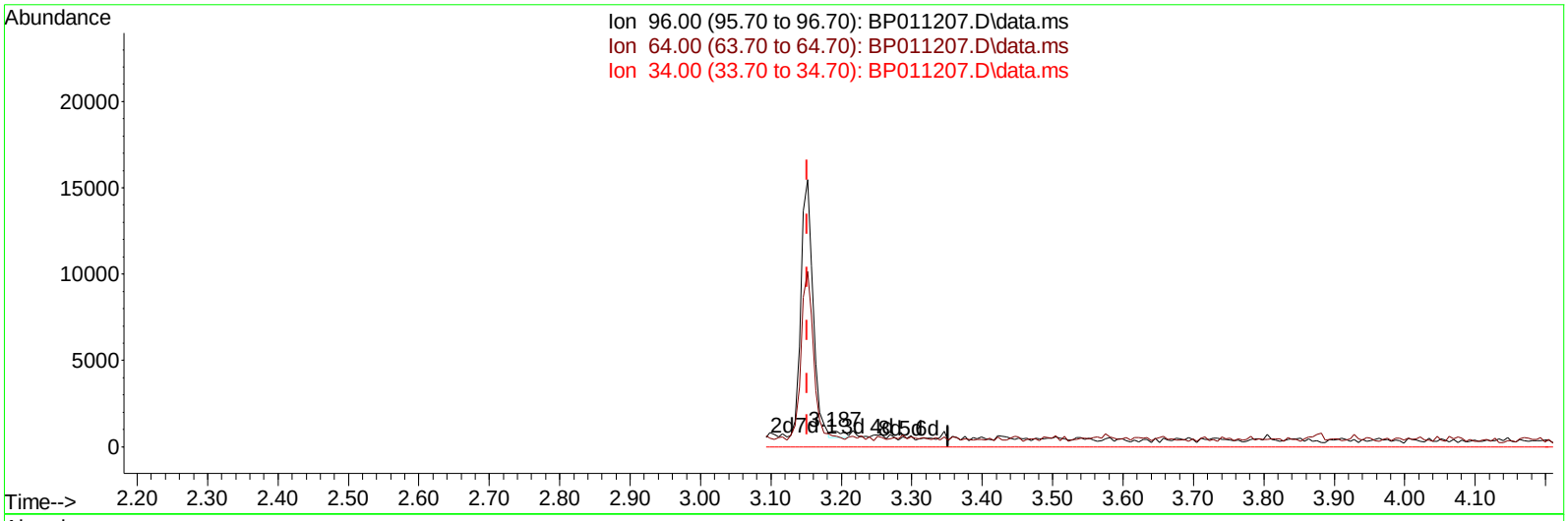
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080122\
Data File : BP011207.D
Acq On : 01 Aug 2022 19:21
Operator : CG/JU
Sample : N3792-10
Misc :
ALS Vial : 16 Sample Multiplier: 1

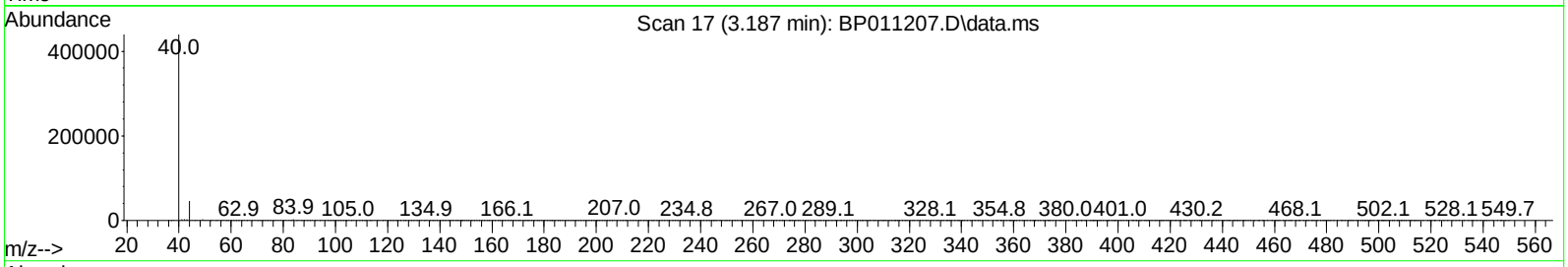
Quant Time: Aug 01 23:01:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP072822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 28 16:05:22 2022
Response via : Initial Calibration

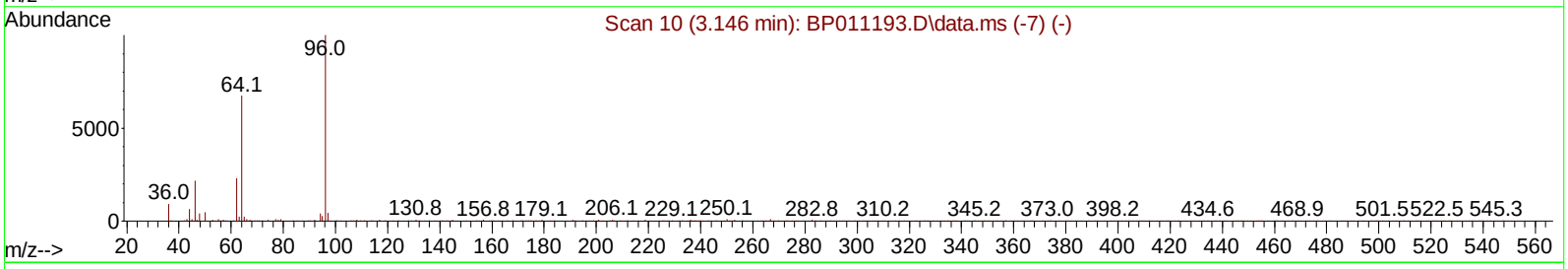
Instrument :
BNA_P
ClientSampleId :
DBUN4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/02/2022
Supervised By :mohammad ahmed 08/03/2022







TIC: BP011207.D\data.ms

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

3.187min (+ 0.035) 0.08 ng/uL

response 500

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	68.31
34.00	0.00	0.00
0.00	0.00	0.00

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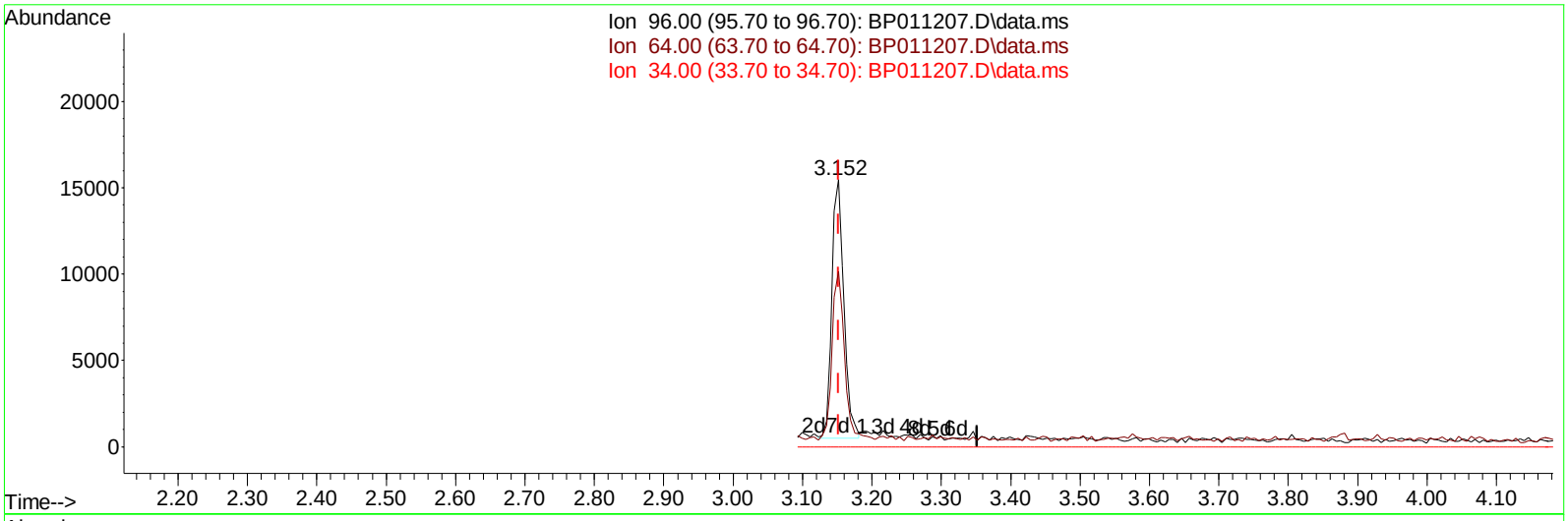

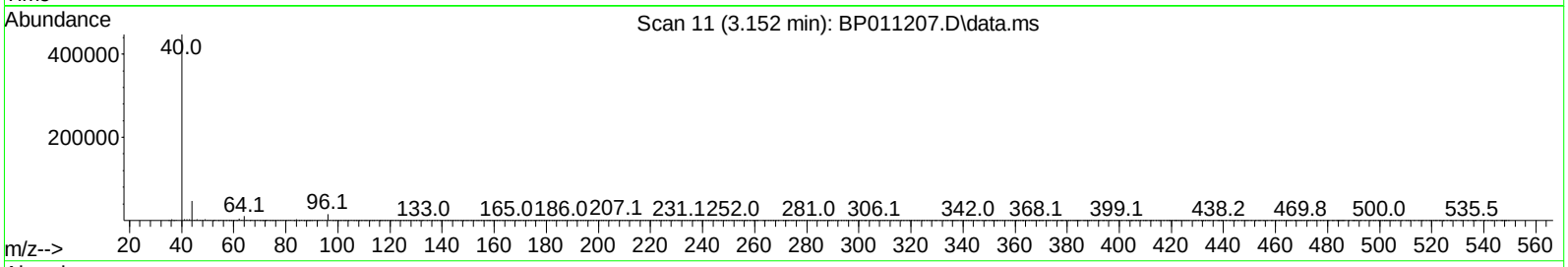

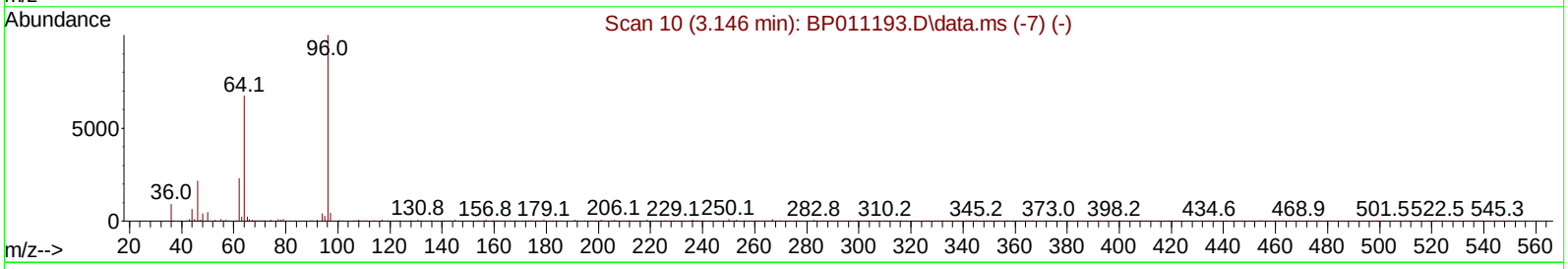

TIC: BP011207.D\data.ms

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

3.152min (-0.000) 2.99 ng/uL m

response 18112

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.90	65.57#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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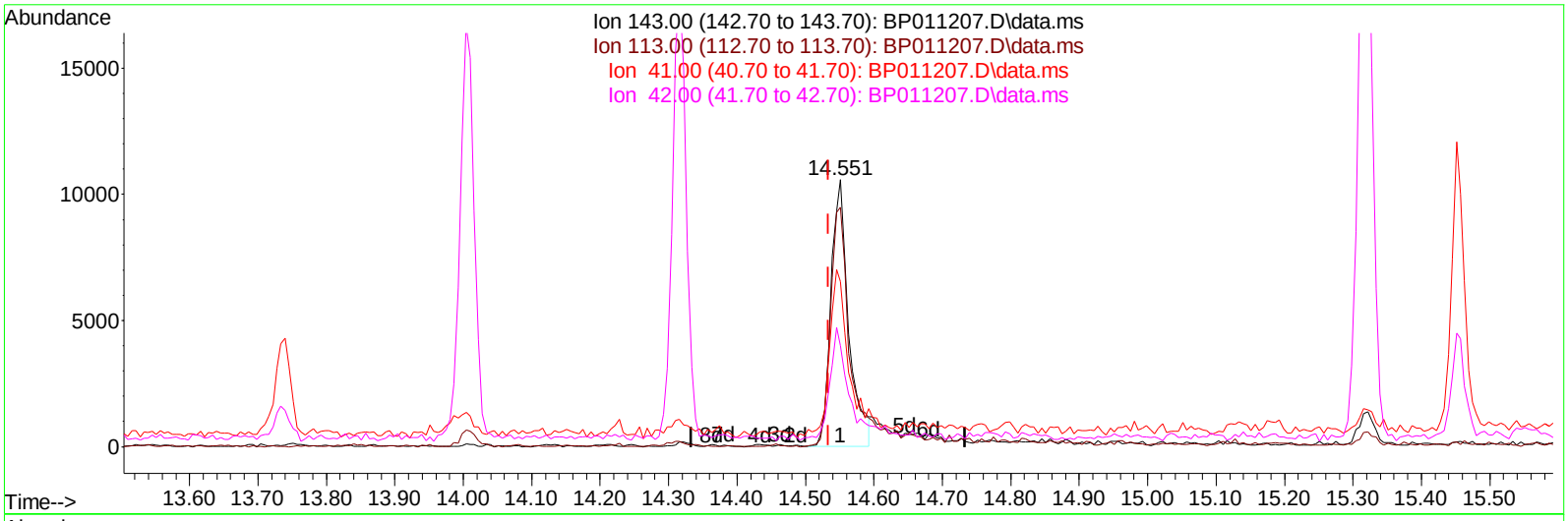

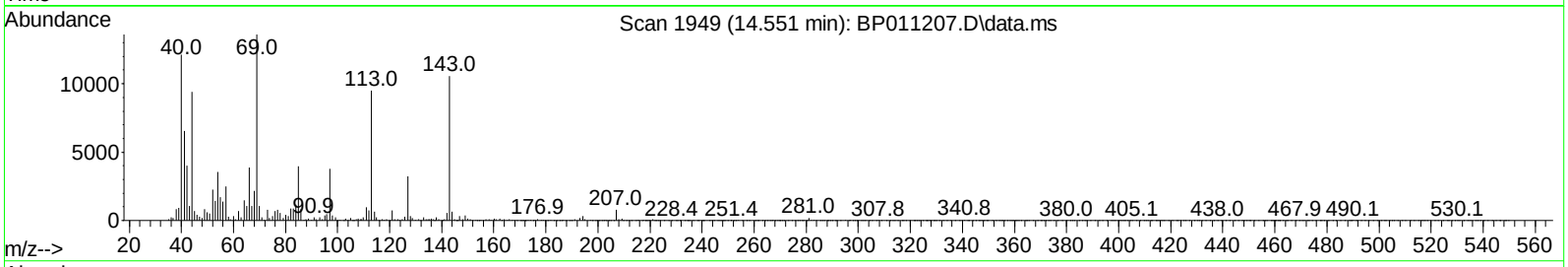

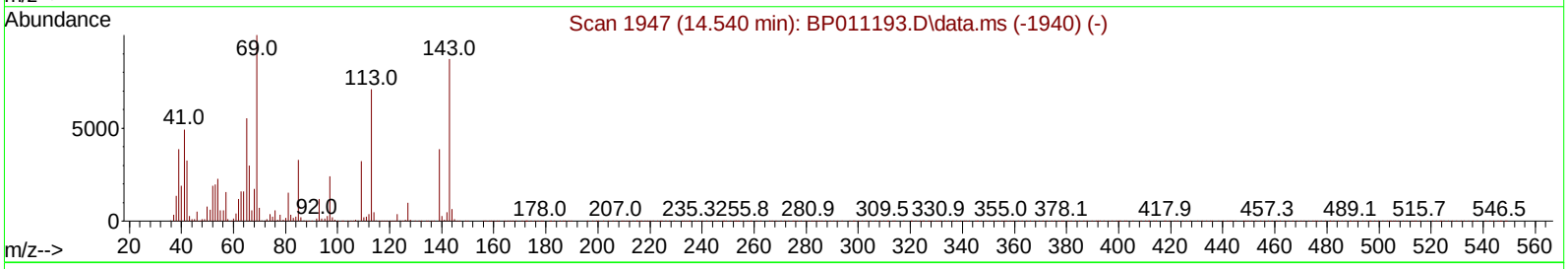

TIC: BP011207.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (+ 0.018) 2.62 ng/ul

response 19007

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	89.79
41.00	59.00	61.83
42.00	41.10	37.76

Quantitation Report (Qedit)

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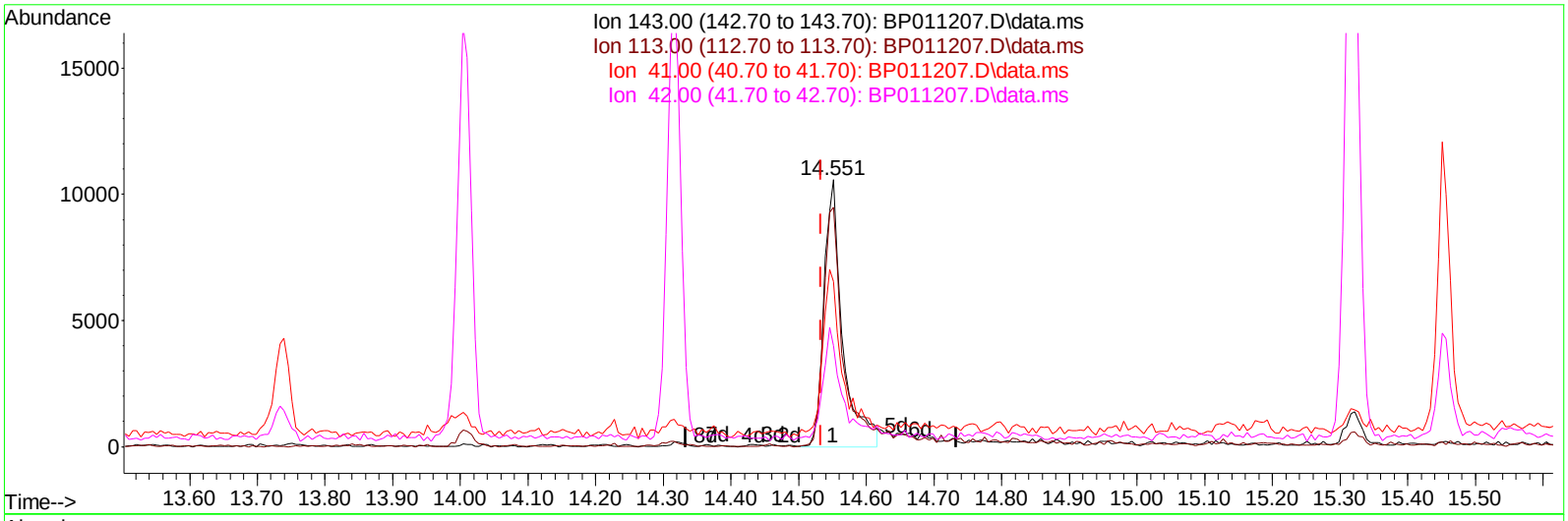

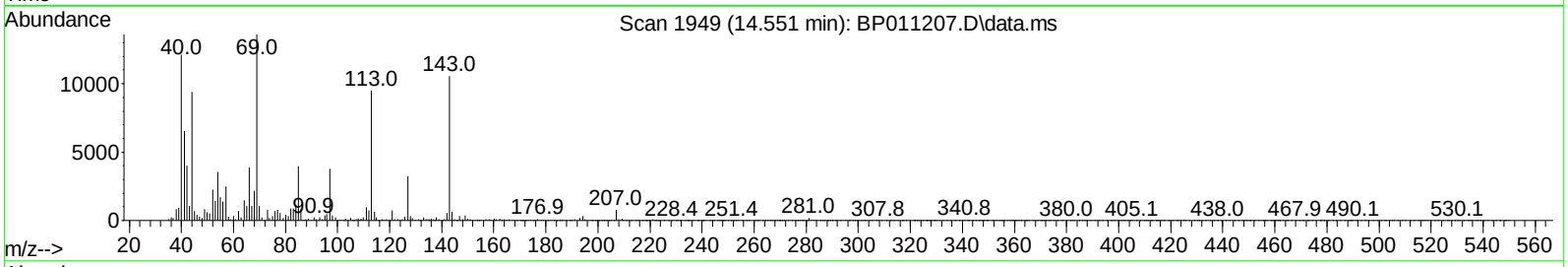

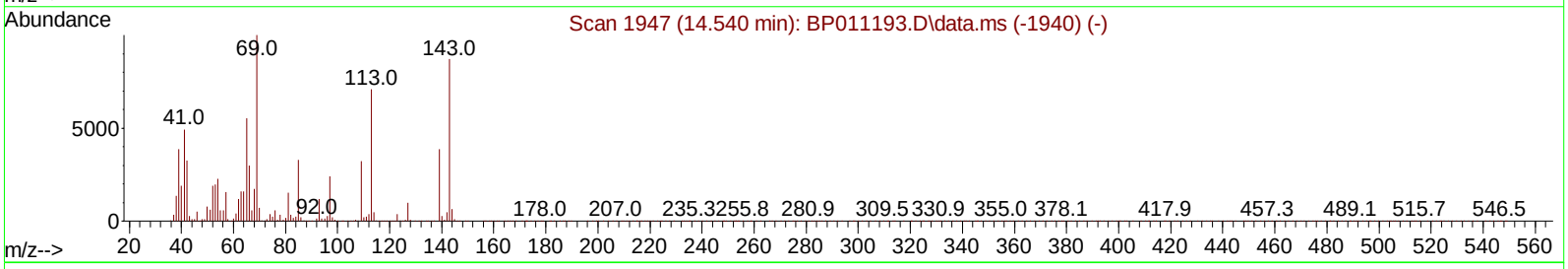

TIC: BP011207.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.551min (+ 0.018) 2.80 ng/ul m

response 20350

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	94.50	89.79
41.00	59.00	61.83
42.00	41.10	37.76

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	277757	20.000	ng/ul	0.00
20) Naphthalene-d8	10.440	136	1142793	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.316	164	592114	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.086	188	1173381	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.210	240	1084204	20.000	ng/ul	# 0.00
88) Perylene-d12	23.562	264	1168558	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	18112m	2.986	ng/uL	0.00
4) Pyridine-d5	3.575	84	37917	2.210	ng/ul	0.02
7) Phenol-d5	6.834	99	123620	5.099	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.993	67	437810	25.539	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	387136	20.263	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	229565	11.575	ng/ul	0.00
21) Nitrobenzene-d5	8.810	128	249861	28.133	ng/ul	0.00
24) 2-Nitrophenol-d4	9.534	143	235040	27.628	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	356819	23.800	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	359136	14.776	ng/ul	0.00
46) Dimethylphthalate-d6	13.739	166	1217874	30.189	ng/ul	0.00
49) Acenaphthylene-d8	14.010	160	1508674	29.760	ng/ul	0.00
54) 4-Nitrophenol-d4	14.551	143	20350m	2.804	ng/ul	0.02
60) Fluorene-d10	15.322	176	1031522	30.429	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	104865	17.729	ng/ul	0.01
73) Anthracene-d10	17.186	188	1653134	31.974	ng/ul	0.00
81) Pyrene-d10	19.445	212	1908173	32.753	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.410	264	1867477	32.175	ng/ul	0.00
Target Compounds						Qvalue

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

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