Quantitation Report (Qedit)

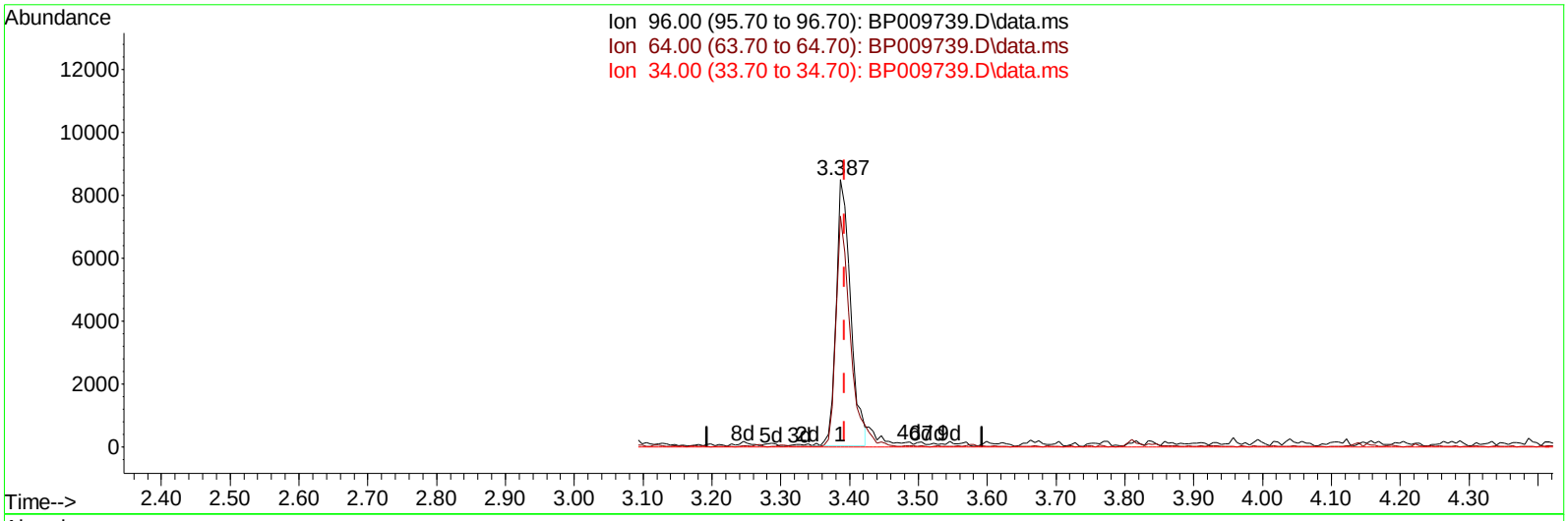
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040822\
Data File : BP009739.D
Acq On : 09 Apr 2022 03:47
Operator : CG/JU
Sample : N2266-05
Misc :
ALS Vial : 34 Sample Multiplier: 1

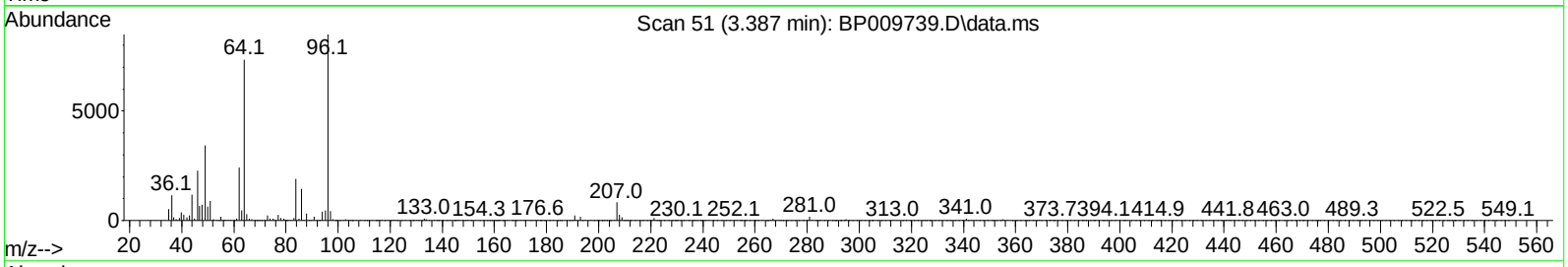
Quant Time: Apr 09 05:53:04 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 07 19:16:52 2022
Response via : Initial Calibration

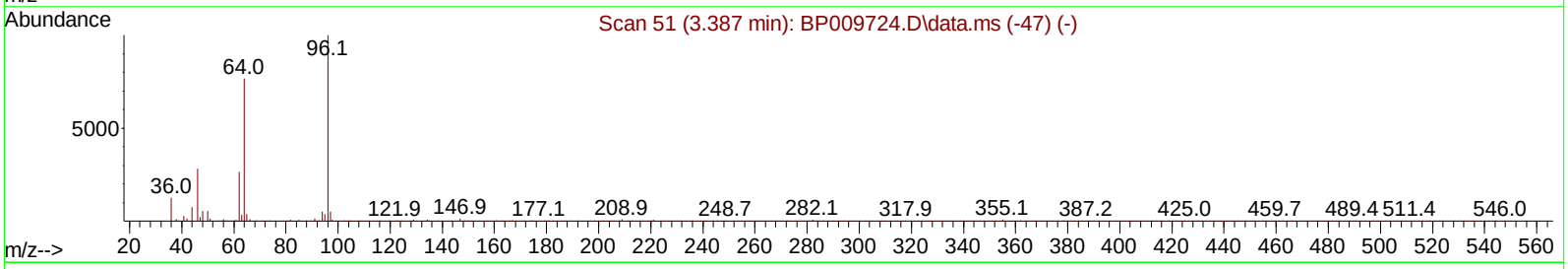
Instrument :
BNA_P
ClientSampleId :
YAZ16

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/11/2022
Supervised By :Yogesh Patel 04/12/2022







TIC: BP009739.D\data.ms

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

3.387min (-0.006) 5.79 ng/uL

response 12288

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.49#
34.00	0.00	0.00
0.00	0.00	0.00

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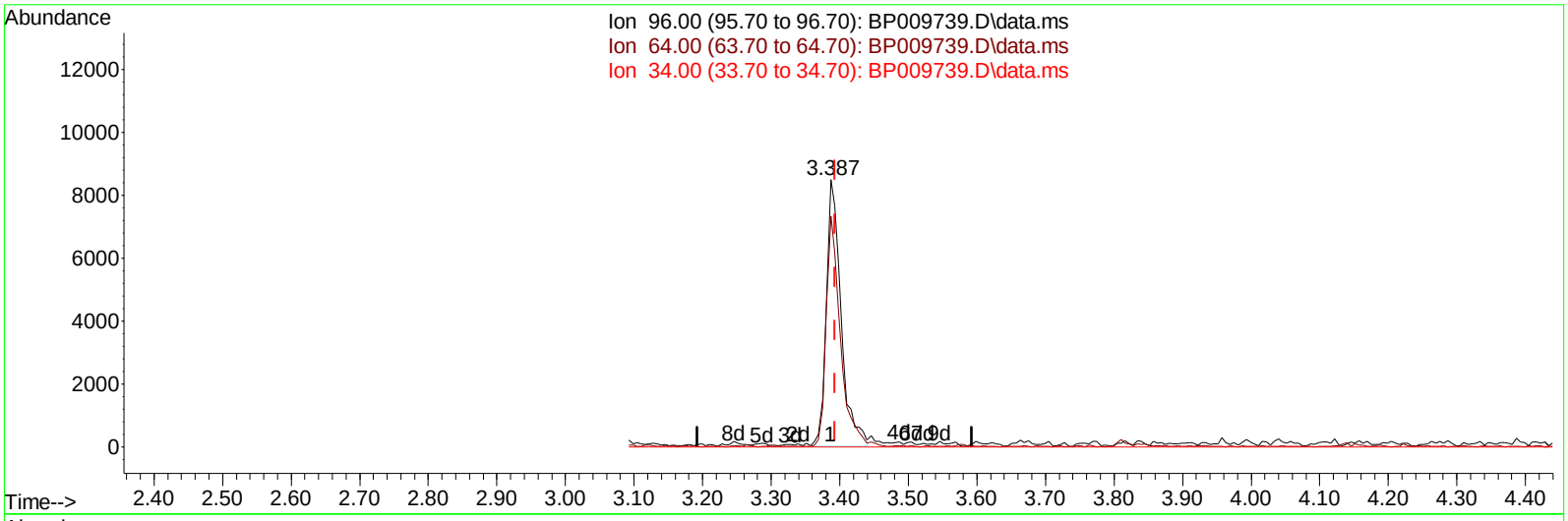

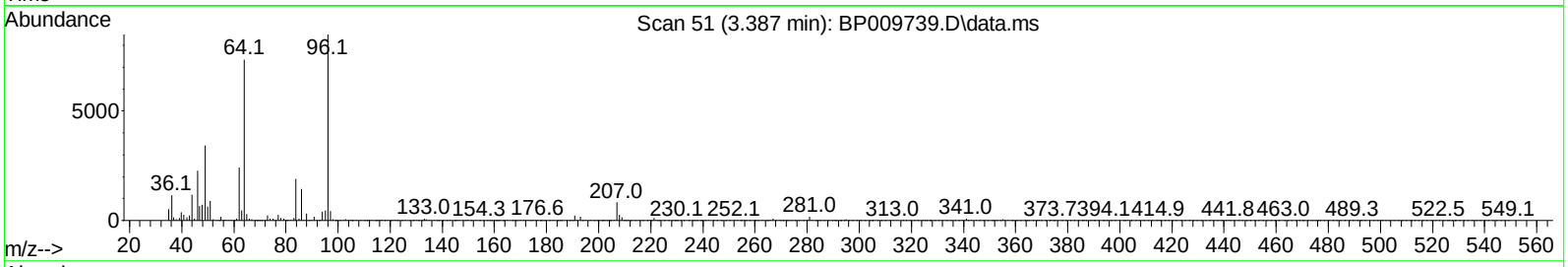

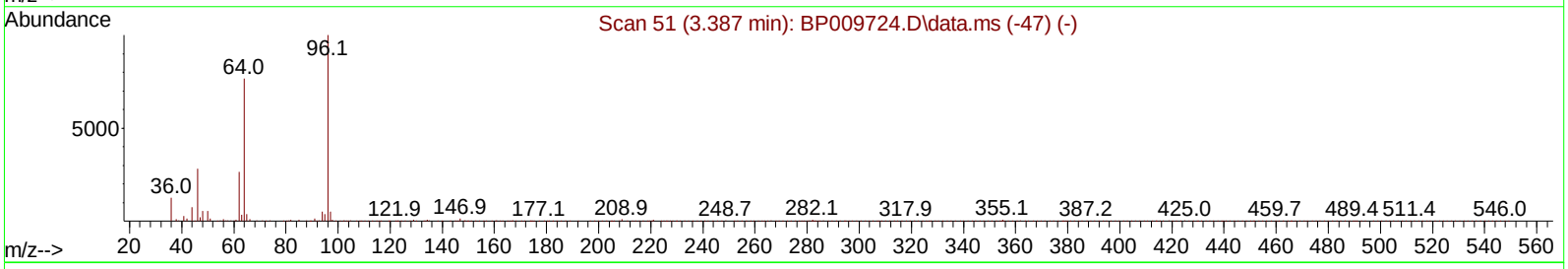

TIC: BP009739.D\data.ms

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

3.387min (-0.006) 6.04 ng/uL m

response 12811

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.49#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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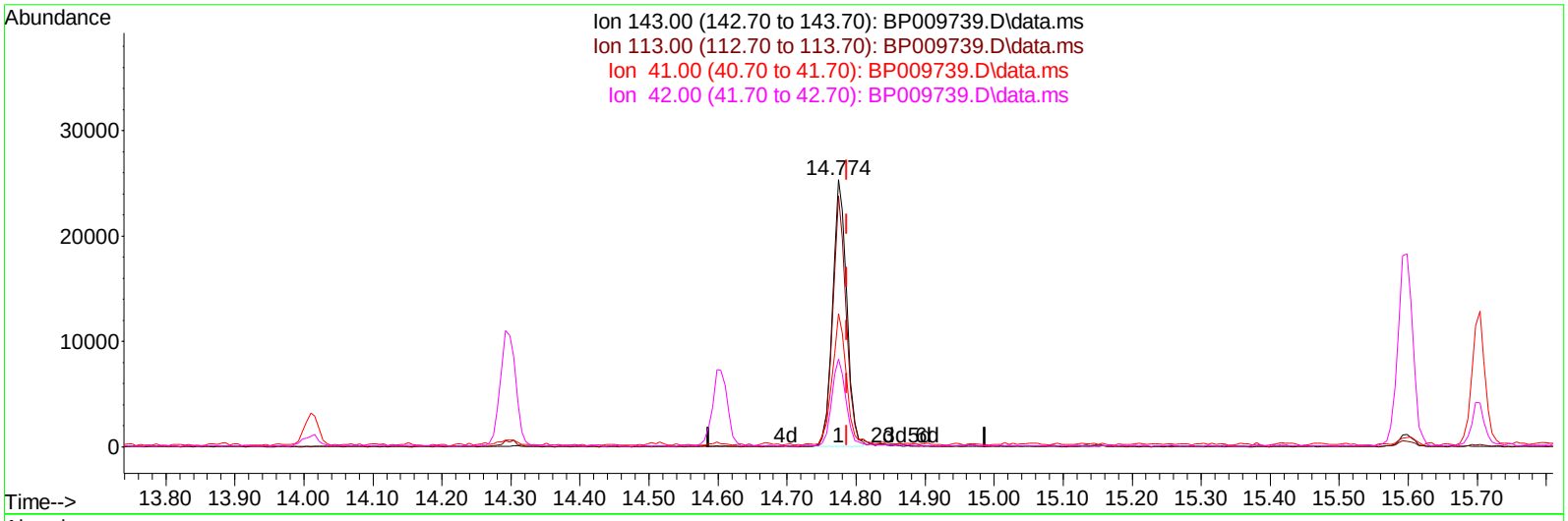

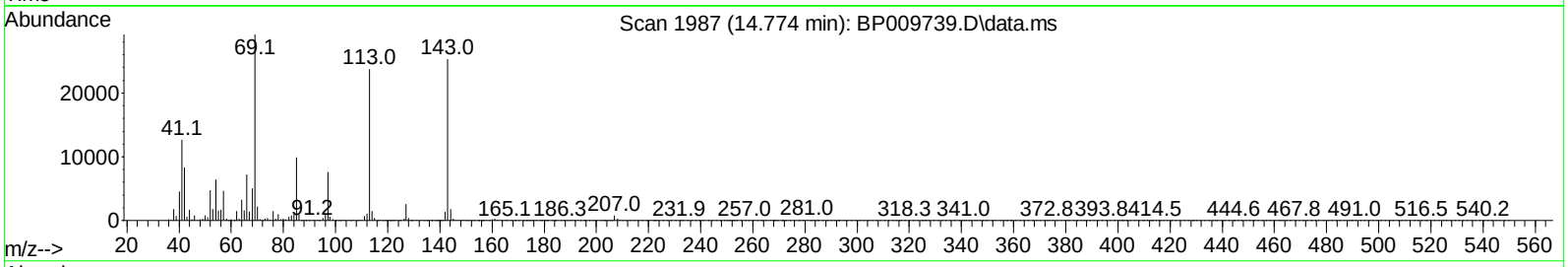

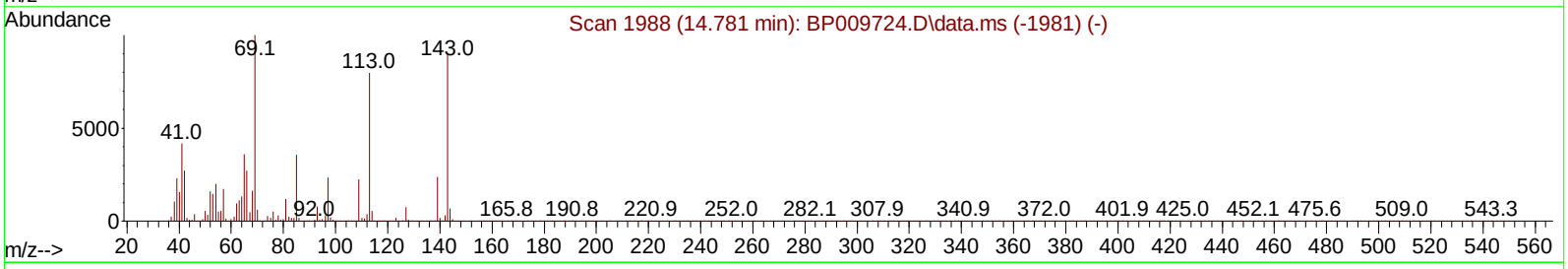

TIC: BP009739.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.774min (-0.012) 8.44 ng/ul

response 36365

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	93.94#
41.00	23.20	49.99#
42.00	18.00	32.81#

Quantitation Report (Qedit)

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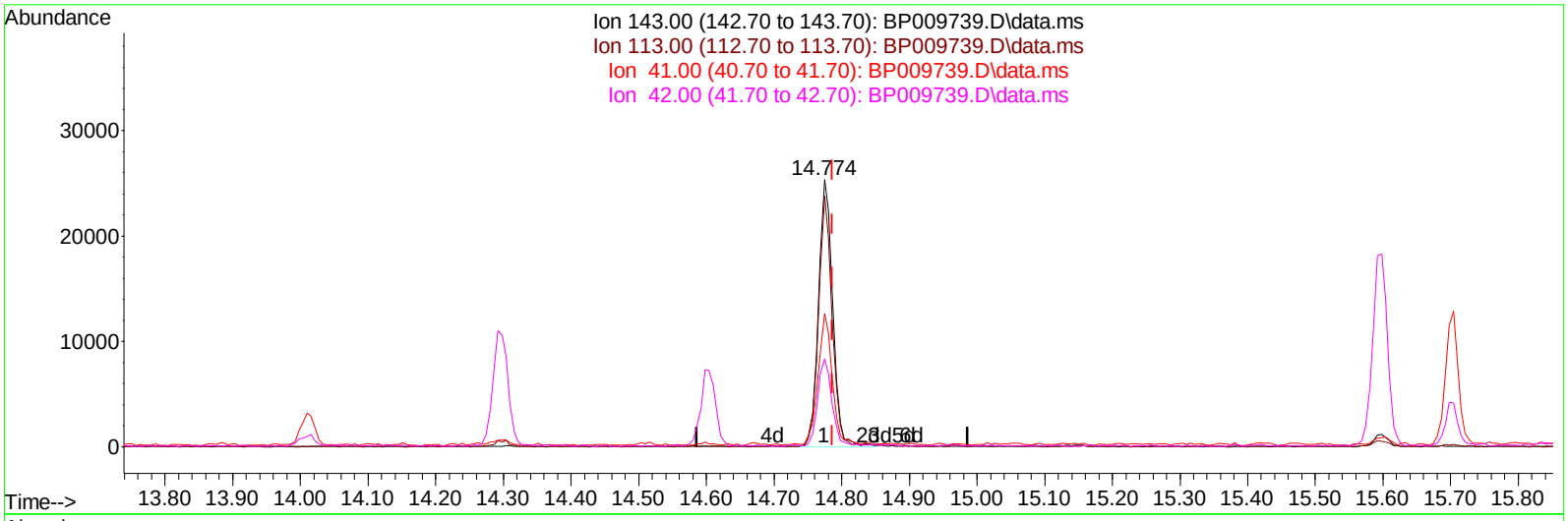

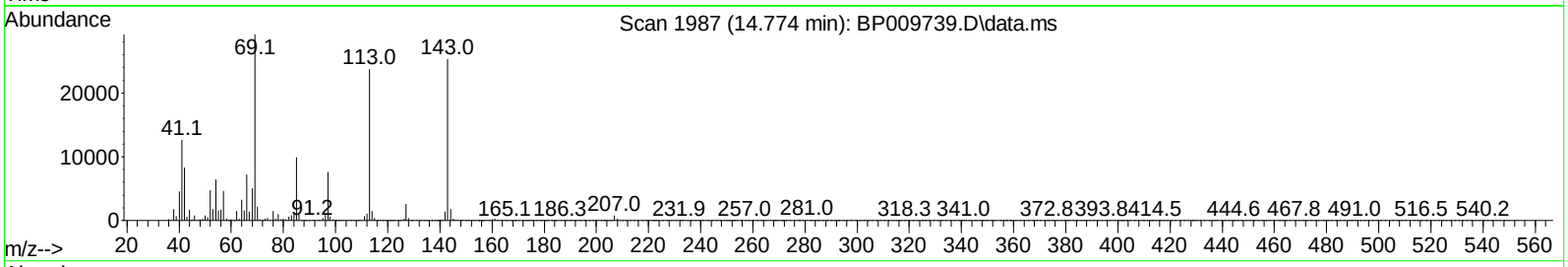

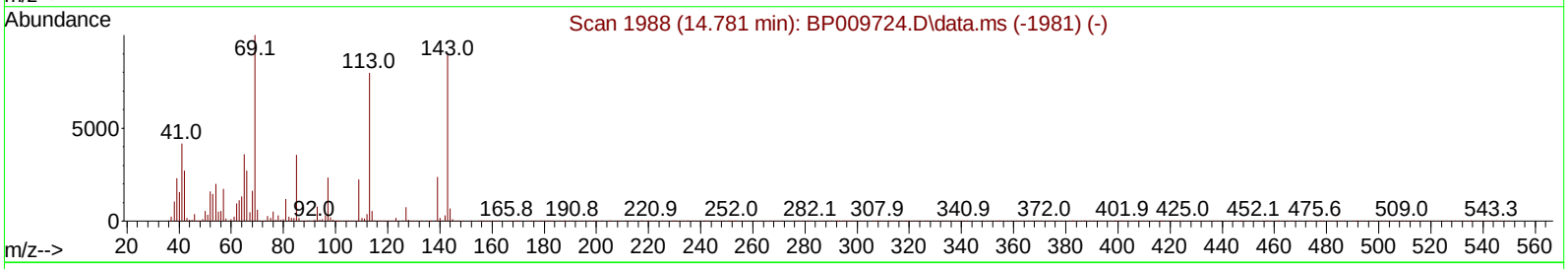

TIC: BP009739.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.774min (-0.012) 8.58 ng/ul m

response 36972

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	93.94#
41.00	23.20	49.99#
42.00	18.00	32.81#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Yogesh Patel 04/12/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	94694	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	422153	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.604	164	302446	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	693119	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.433	240	746980	20.000	ng/ul	-0.01
88) Perylene-d12	23.909	264	784536	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	12811m	6.038	ng/uL	0.00
4) Pyridine-d5	3.804	84	67306	11.505	ng/ul	0.00
7) Phenol-d5	7.122	99	71207	8.565	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.298	67	224565	45.373	ng/ul	0.00
11) 2-Chlorophenol-d4	7.498	132	212702	33.918	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	155737	22.440	ng/ul	-0.01
21) Nitrobenzene-d5	9.128	128	148477	48.773	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.857	143	156249	46.815	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	266152	38.108	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	371936	37.390	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1148574	48.346	ng/ul	-0.01
49) Acenaphthylene-d8	14.298	160	1237567	43.970	ng/ul	0.00
54) 4-Nitrophenol-d4	14.774	143	36972m	8.581	ng/ul	-0.01
60) Fluorene-d10	15.598	176	957583	47.913	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	173961	42.367	ng/ul	0.00
73) Anthracene-d10	17.457	188	1619629	48.805	ng/ul	0.00
81) Pyrene-d10	19.680	212	1986362	48.689	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.750	264	2046700	50.096	ng/ul	-0.01
Target Compounds						Qvalue

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

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