Quantitation Report (Qedit)

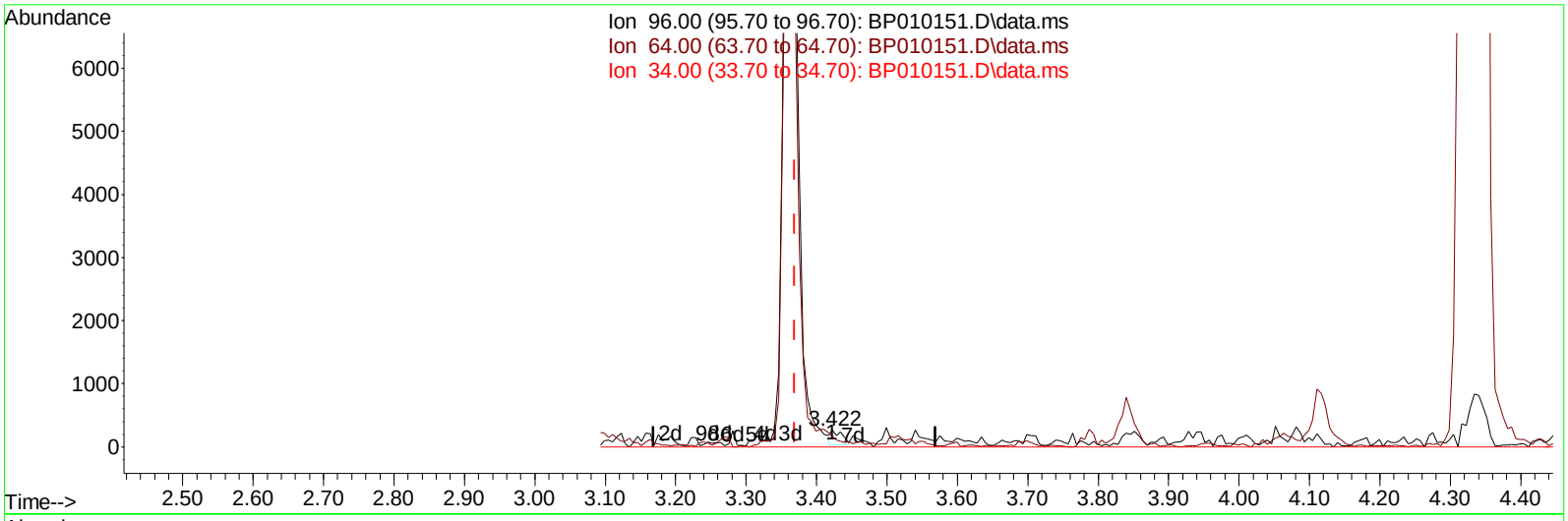
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010151.D
Acq On : 29 Apr 2022 15:02
Operator : CG/JU
Sample : N2587-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

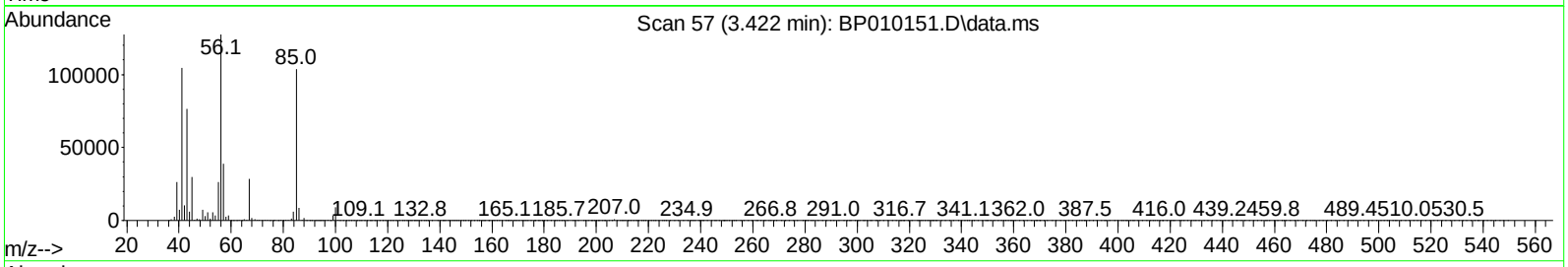
Quant Time: Apr 30 00:30:29 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

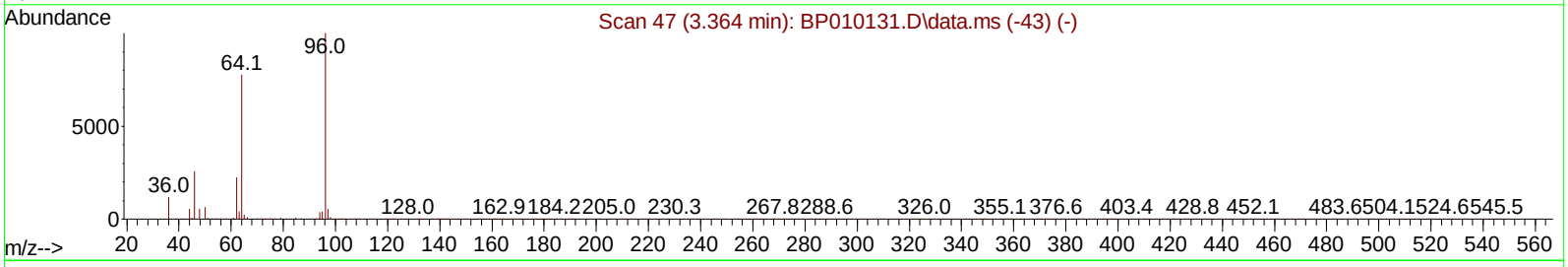
Instrument :
BNA_P
ClientSampleId :
EXET3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/02/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010151.D\data.ms

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

3.422min (+ 0.053) 0.07 ng/uL

response 238

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	61.60
34.00	0.00	0.00
0.00	0.00	0.00

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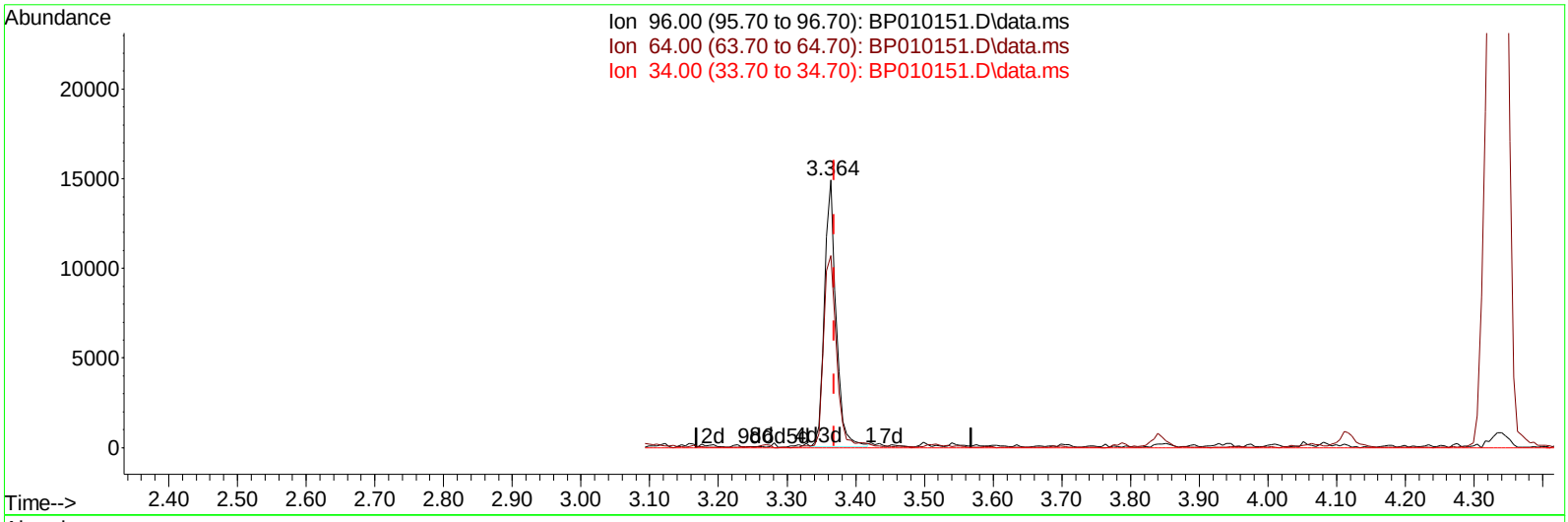

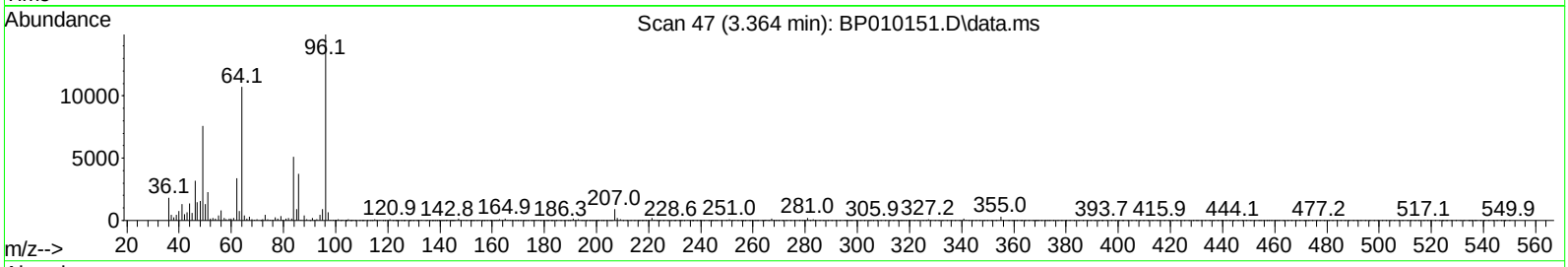

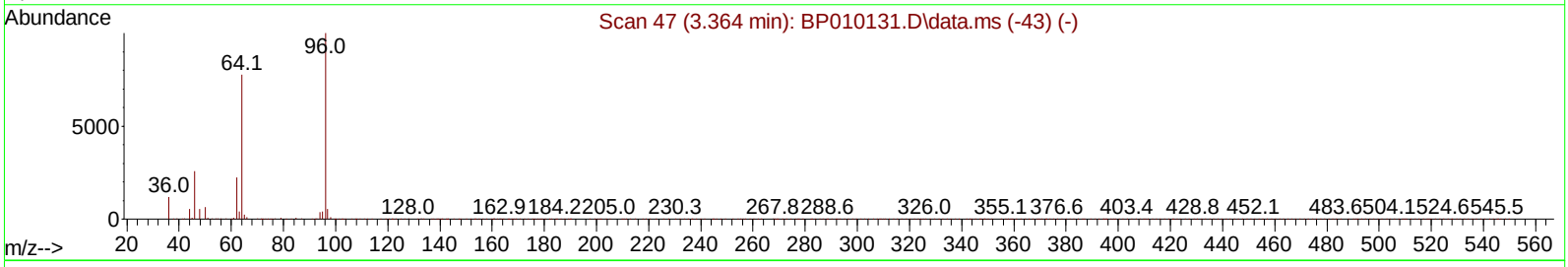

TIC: BP010151.D\data.ms

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

3.364min (-0.005) 5.35 ng/uL m

response 17404

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	71.82#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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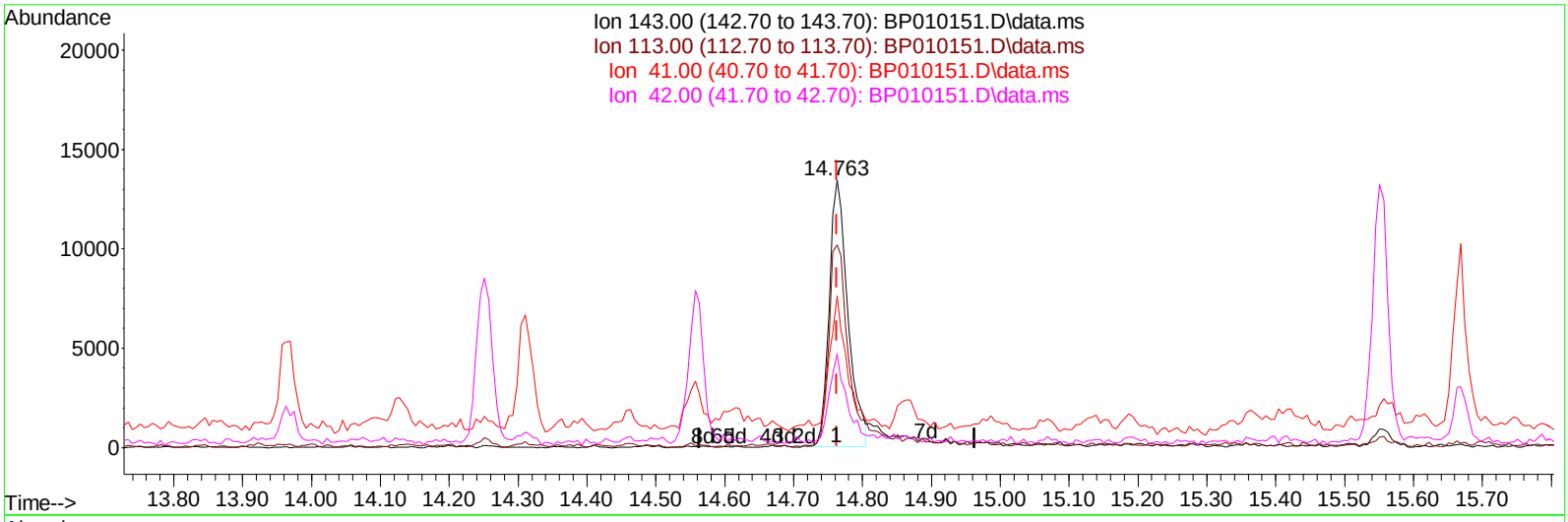

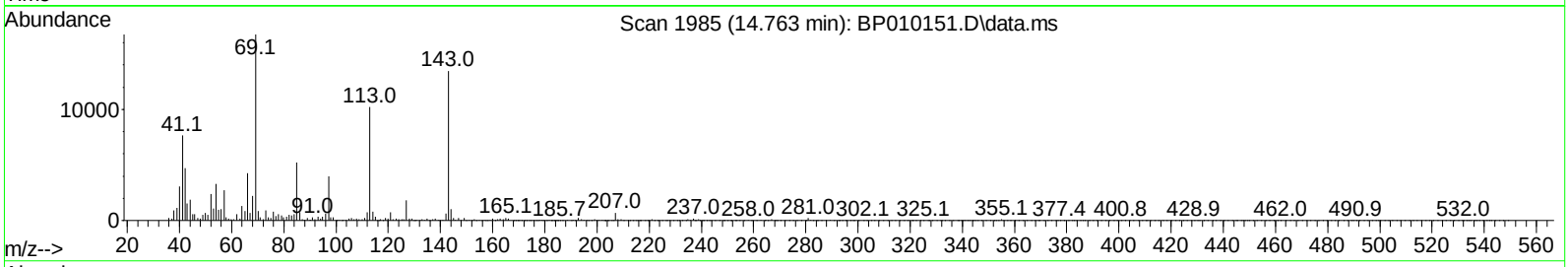

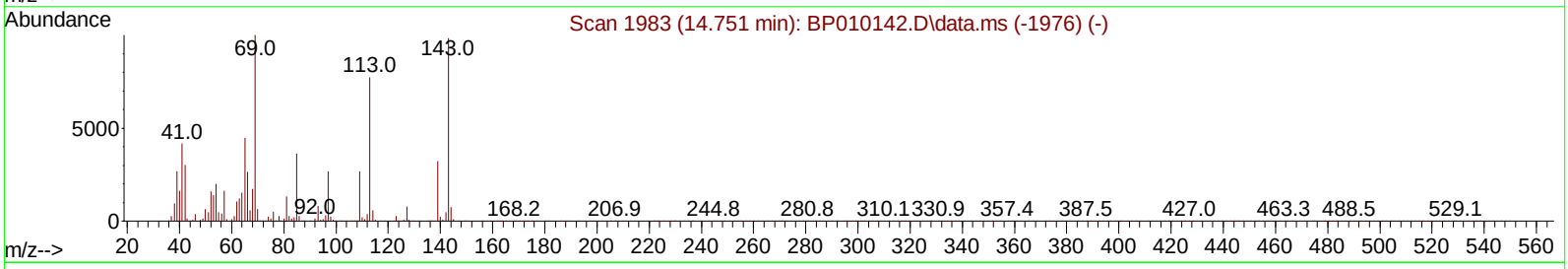

TIC: BP010151.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 5.14 ng/ul

response 25046

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.86#
41.00	23.20	56.81#
42.00	18.00	35.08#

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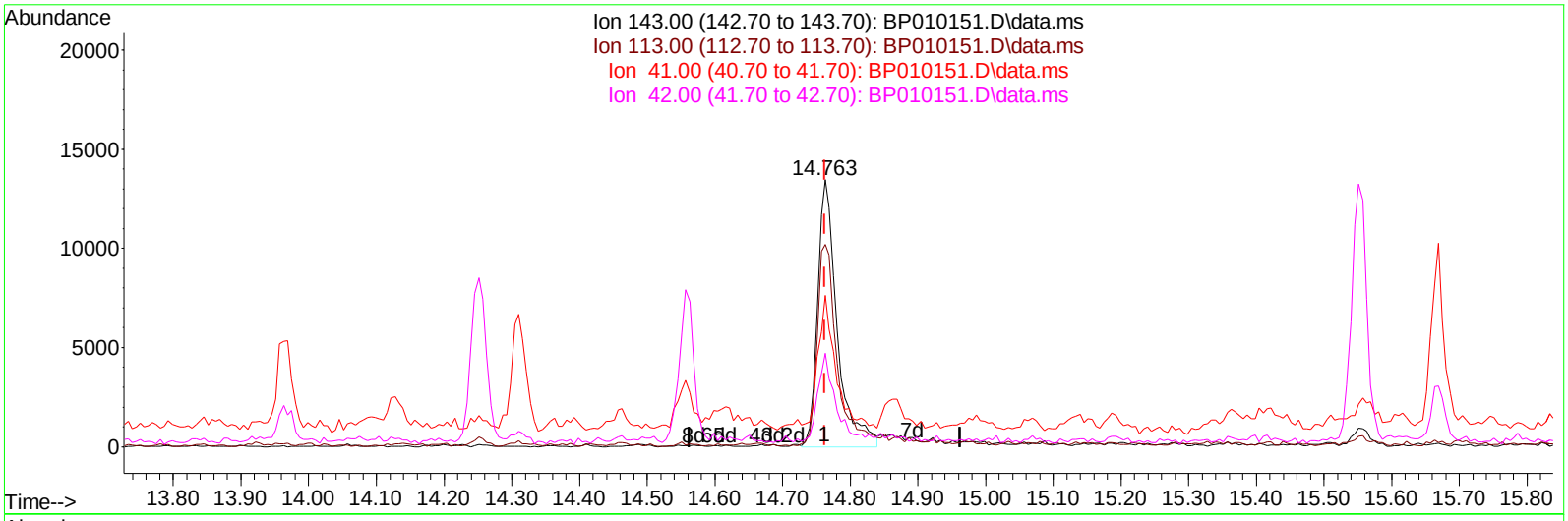

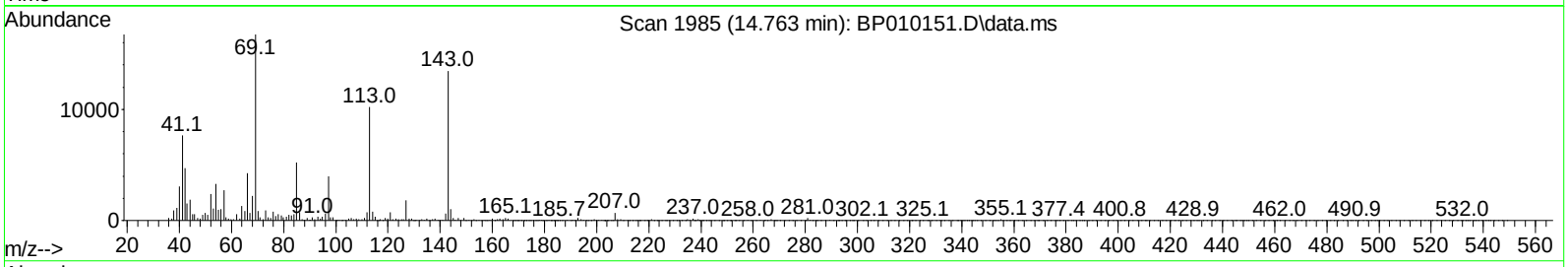

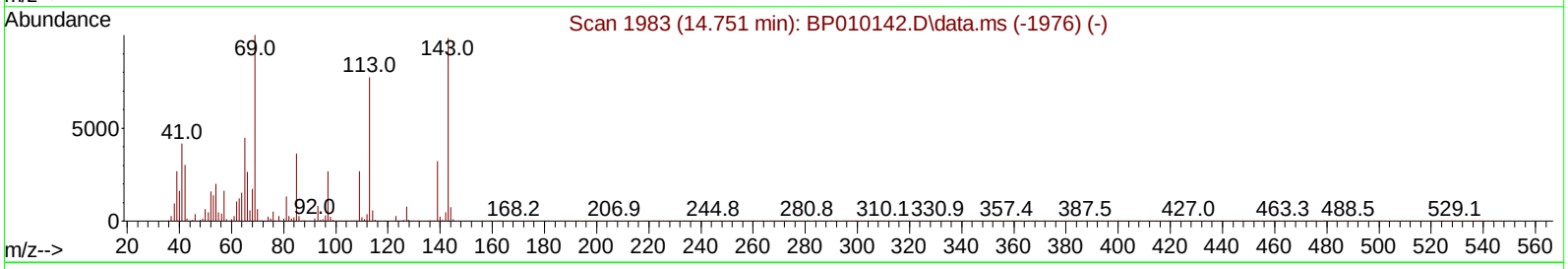

TIC: BP010151.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 5.57 ng/ul m

response 27115

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	75.86#
41.00	23.20	56.81#
42.00	18.00	35.08#

Quantitation Report (QT Reviewed)

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Reviewed By :Jagrut Upadhyay 05/02/2022
Supervised By :Yogesh Patel 05/05/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	141824	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.728	136	559099	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.557	164	340021	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.310	188	656167	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.398	240	619083	20.000	ng/ul	#-0.01
88) Perylene-d12	23.845	264	627494	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	17404m	5.351	ng/uL	0.00
4) Pyridine-d5	3.781	84	100870	11.025	ng/ul	0.00
7) Phenol-d5	7.093	99	99516	7.714	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	246996	31.909	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.457	132	247181	25.873	ng/ul	-0.01
15) 4-Methylphenol-d8	8.634	113	190626	17.835	ng/ul	-0.02
21) Nitrobenzene-d5	9.087	128	150097	33.754	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.810	143	161978	33.353	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	267468	28.290	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	321594	23.723	ng/ul	-0.02
46) Dimethylphthalate-d6	13.969	166	917564	34.691	ng/ul	-0.01
49) Acenaphthylene-d8	14.251	160	1036326	32.374	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.763	143	27115m	5.569	ng/ul	0.00
60) Fluorene-d10	15.551	176	776083	34.110	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.669	200	130178	31.375	ng/ul	-0.01
73) Anthracene-d10	17.410	188	1170373	36.868	ng/ul	-0.01
81) Pyrene-d10	19.645	212	1322164	37.554	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1240803	37.718	ng/ul	-0.01
Target Compounds						Qvalue

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

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