Quantitation Report (Qedit)

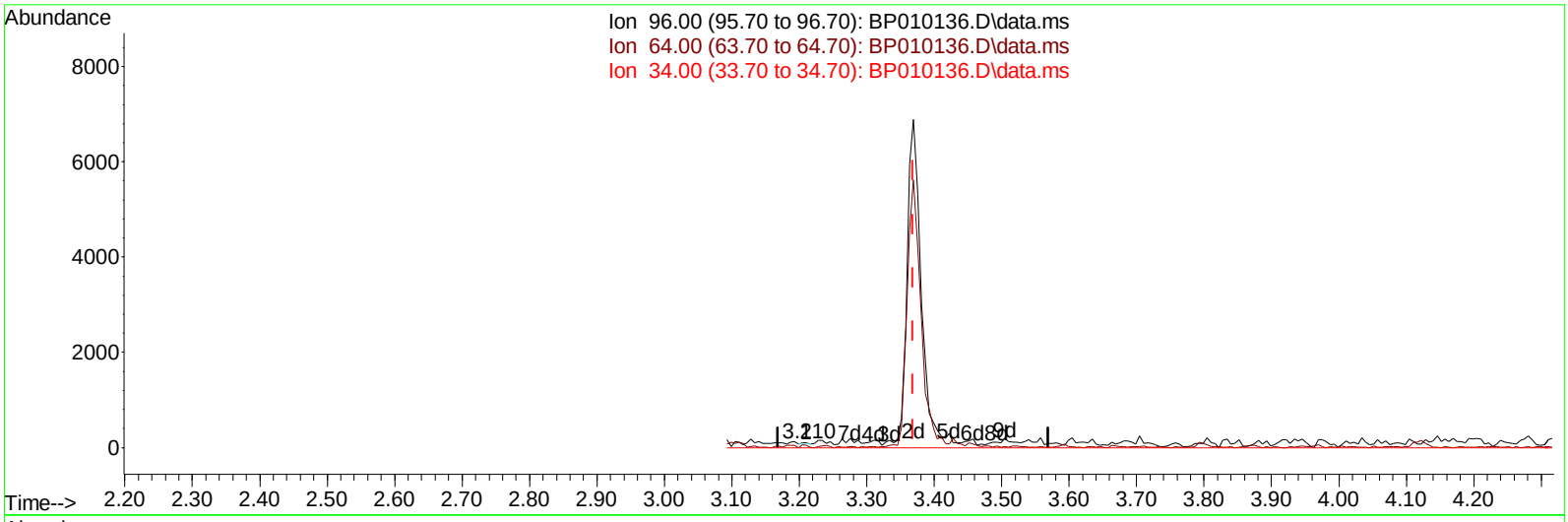
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042822\
Data File : BP010136.D
Acq On : 28 Apr 2022 12:15
Operator : CG/JU
Sample : N2469-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

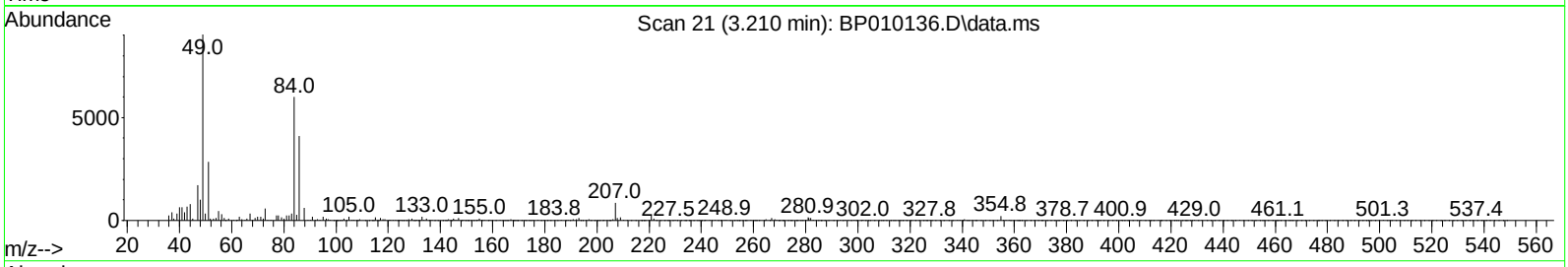
Quant Time: Apr 28 14:08:04 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

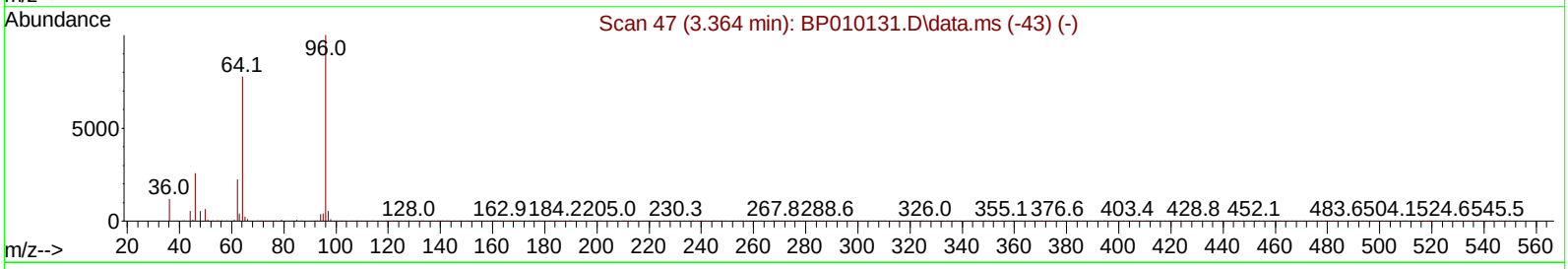
Instrument :
BNA_P
ClientSampleId :
BFFG7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/29/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010136.D\data.ms

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

3.210min (-0.159) 0.00 ng/uL

response 16

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	51.04
34.00	0.00	0.00
0.00	0.00	0.00

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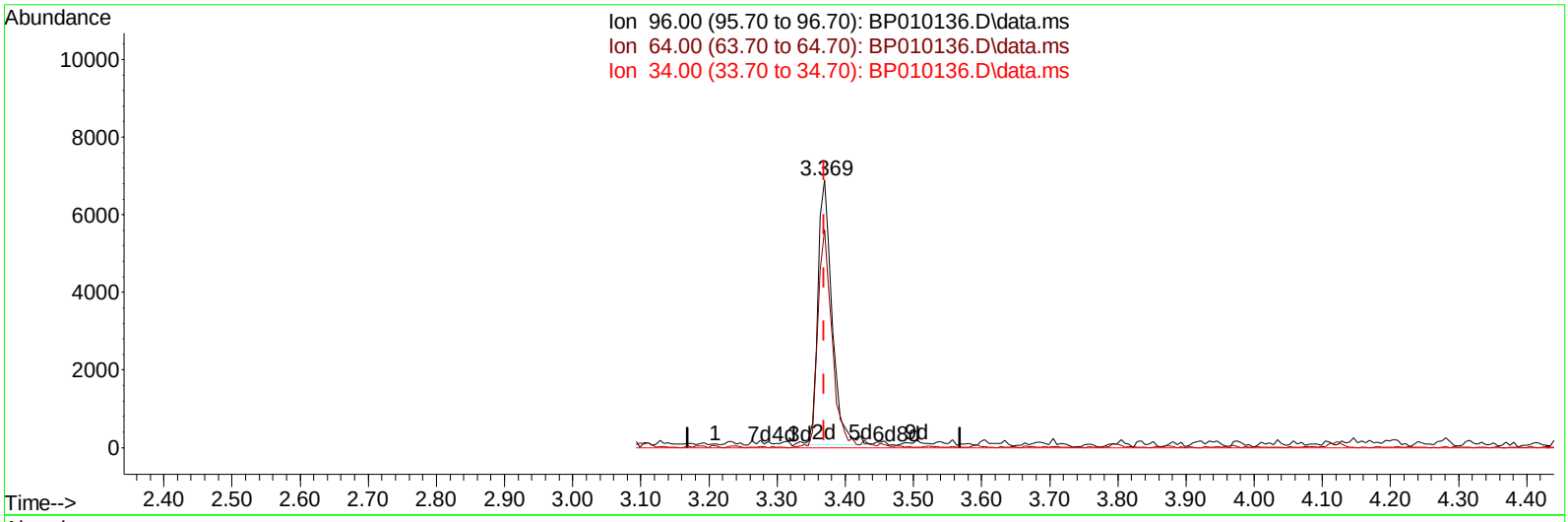

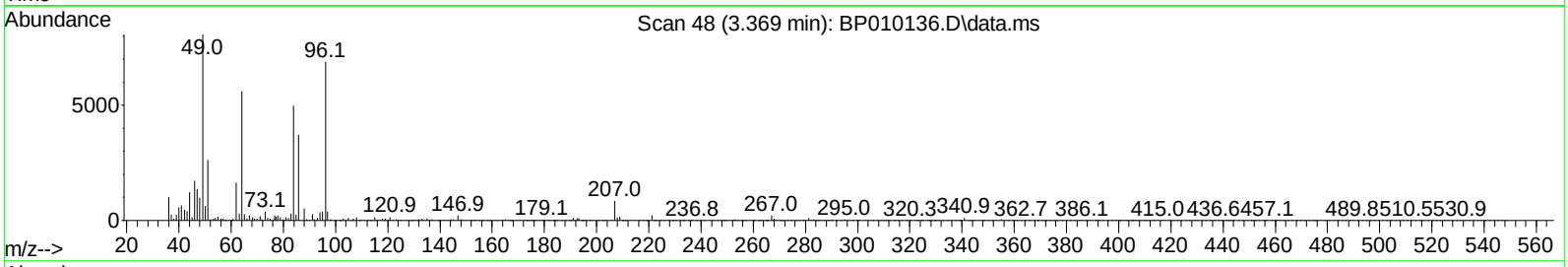

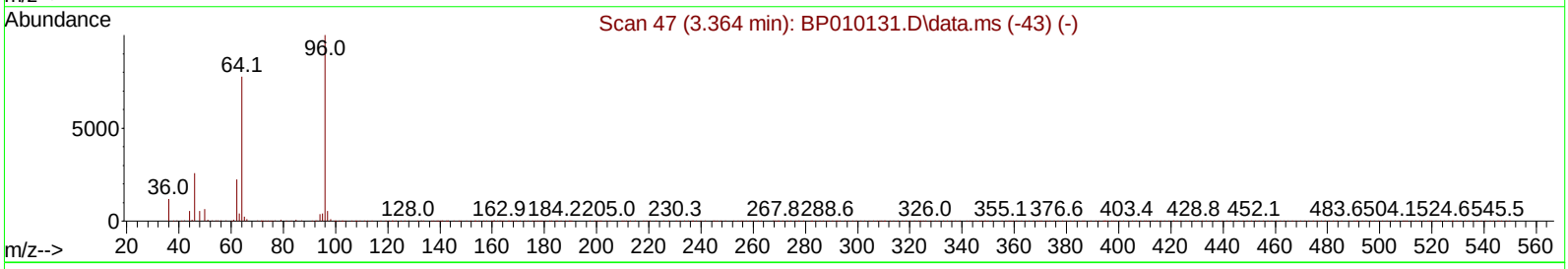

TIC: BP010136.D\data.ms

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

3.369min (+ 0.000) 2.21 ng/uL m

response 9743

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	81.49#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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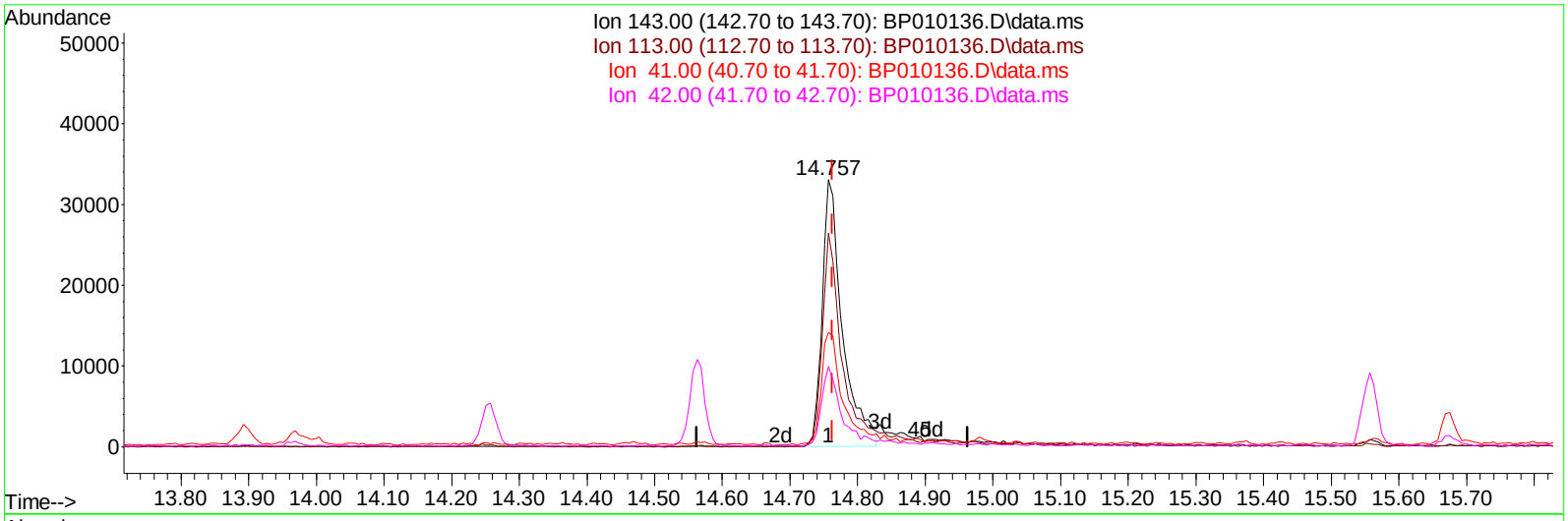

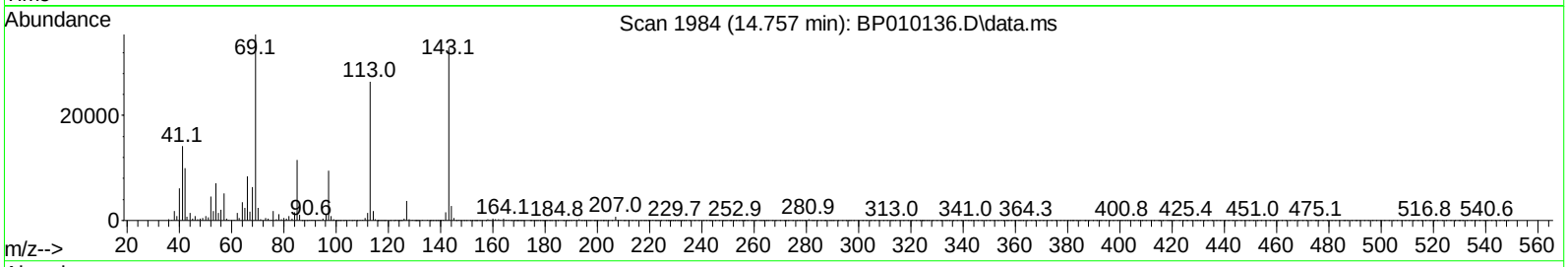

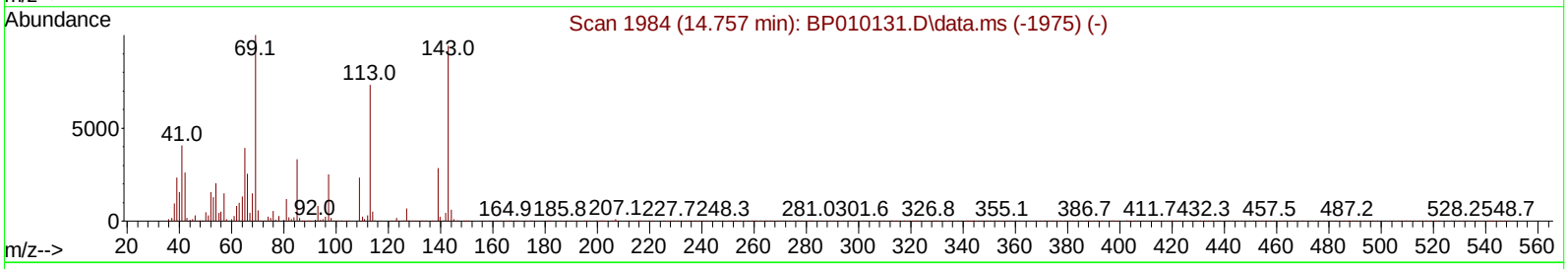

TIC: BP010136.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 8.71 ng/ul

response 69200

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.94#
41.00	23.20	42.85#
42.00	18.00	30.05#

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ClientSampleId :
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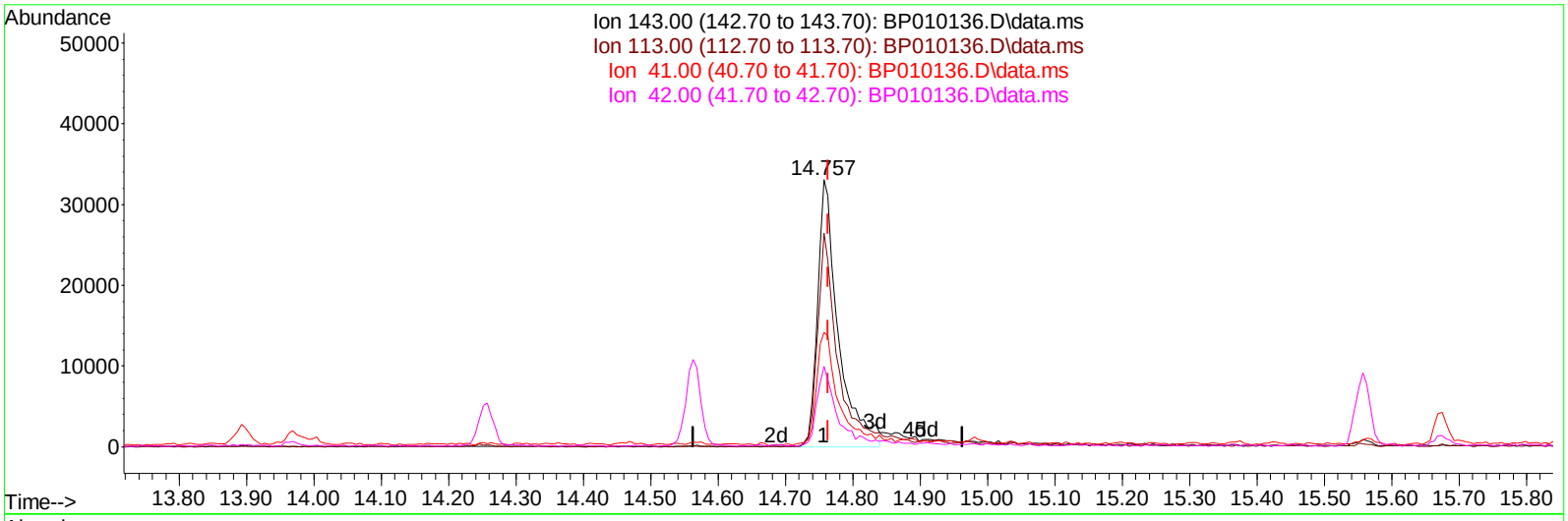

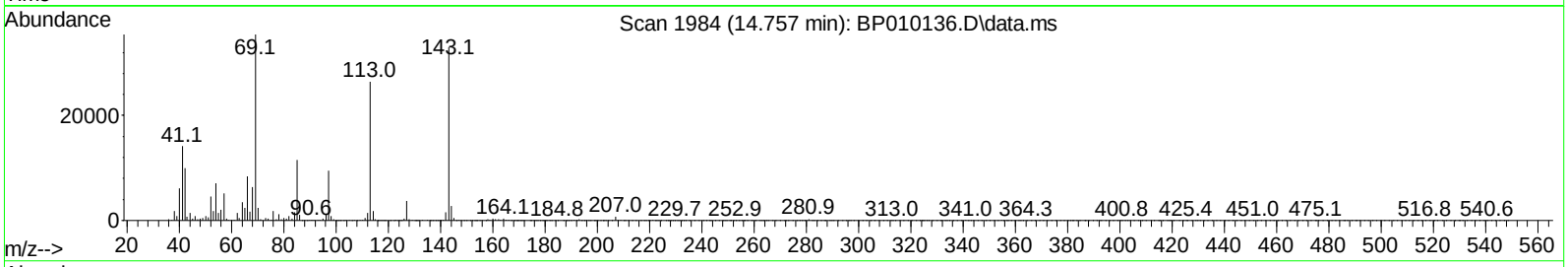

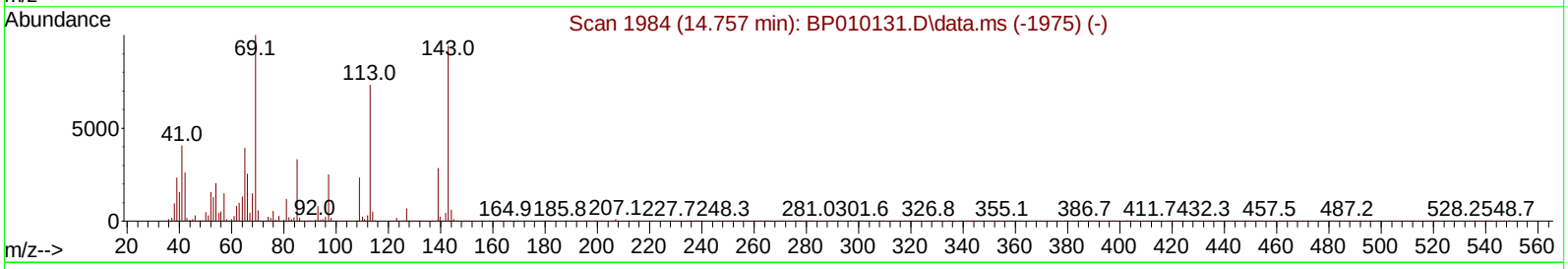

TIC: BP010136.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 8.94 ng/ul m

response 71000

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	79.94#
41.00	23.20	42.85#
42.00	18.00	30.05#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	191894	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	834467	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.563	164	554698	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	1096617	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	713620	20.000	ng/ul	# 0.00
88) Perylene-d12	23.856	264	595030	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	9743m	2.214	ng/uL	0.00
4) Pyridine-d5	3.793	84	43751	3.534	ng/ul	0.00
7) Phenol-d5	7.093	99	196115	11.236	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.257	67	133950	12.789	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	163043	12.613	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	166007	11.479	ng/ul	-0.02
21) Nitrobenzene-d5	9.093	128	85562	12.892	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	92773	12.799	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	162895	11.544	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	186976	9.241	ng/ul	-0.01
46) Dimethylphthalate-d6	13.969	166	629480	14.588	ng/ul	-0.01
49) Acenaphthylene-d8	14.257	160	679887	13.019	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	71000m	8.938	ng/ul	0.00
60) Fluorene-d10	15.557	176	531275	14.313	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	62374	8.995	ng/ul	0.00
73) Anthracene-d10	17.416	188	843062	15.891	ng/ul	0.00
81) Pyrene-d10	19.651	212	962327	23.712	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	572946	18.367	ng/ul	0.00
Target Compounds						Qvalue
82) Pyrene	19.674	202	48906	1.037	ng/ul	99

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

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