Quantitation Report (Qedit)

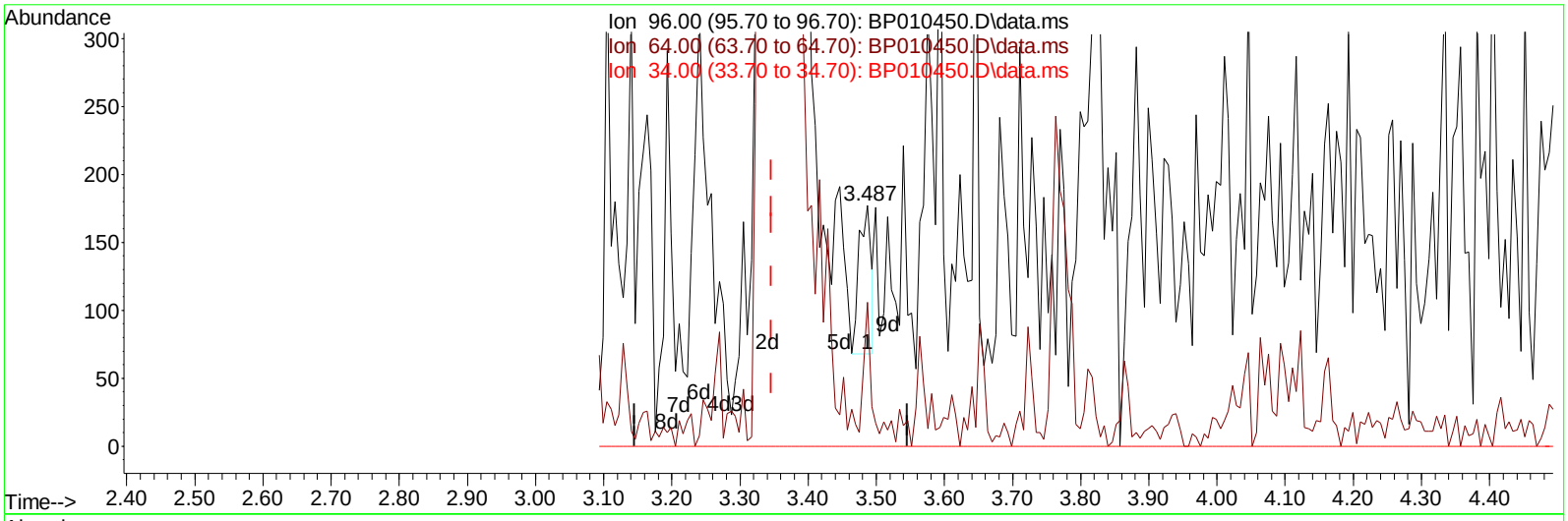
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
Data File : BP010450.D
Acq On : 20 May 2022 23:40
Operator : CG/JU
Sample : N2843-19
Misc :
ALS Vial : 25 Sample Multiplier: 1

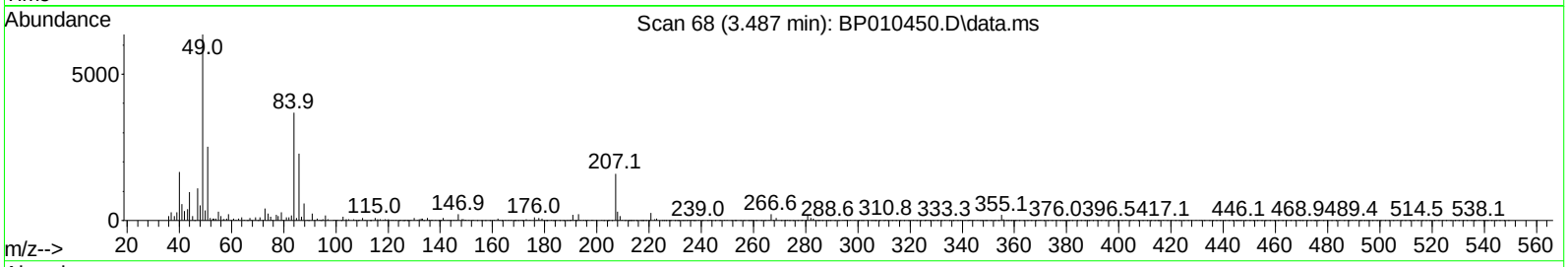
Quant Time: May 21 01:20:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

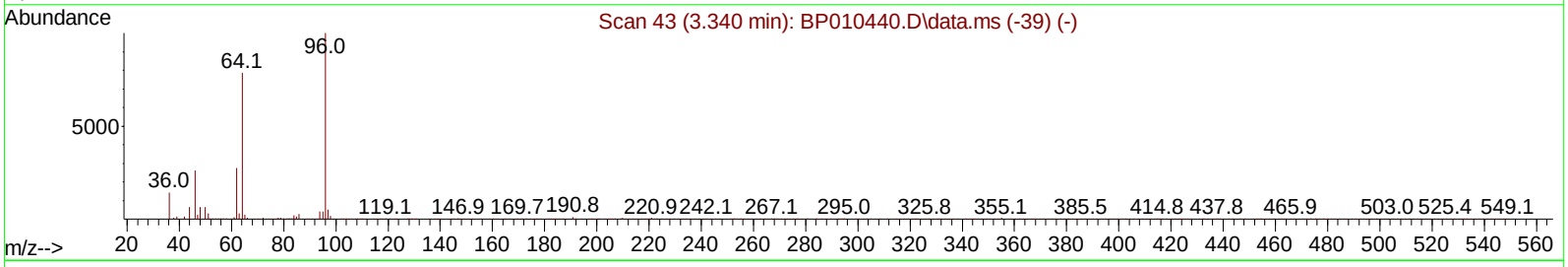
Instrument :
BNA_P
ClientSampleId :
EW4J9

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/23/2022
Supervised By :mohammad ahmed 05/25/2022







TIC: BP010450.D\data.ms

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

3.487min (+ 0.141) 0.04 ng/uL

response 132

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.89
34.00	0.00	0.00
0.00	0.00	0.00

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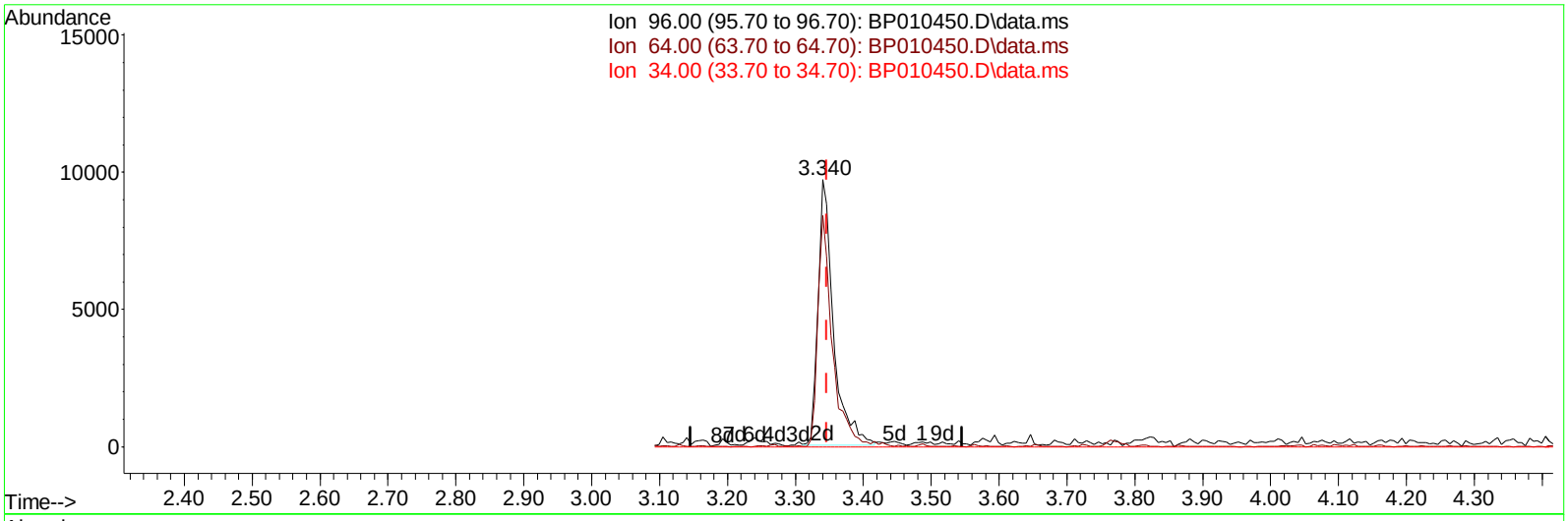

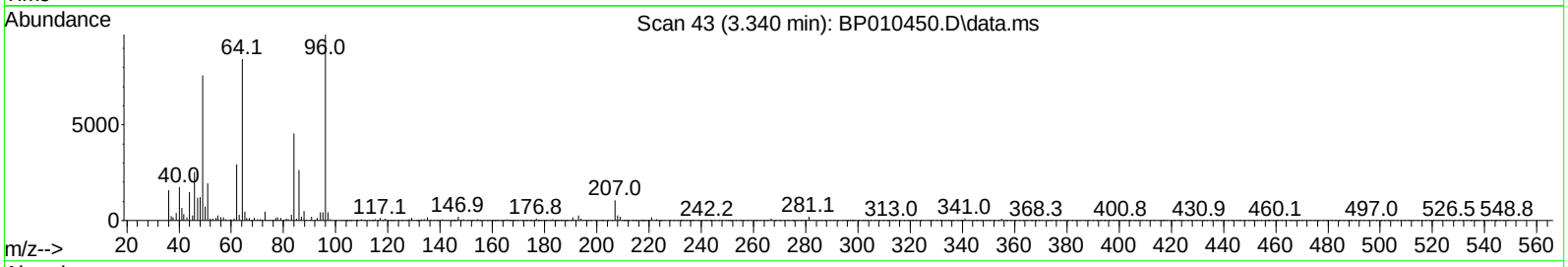

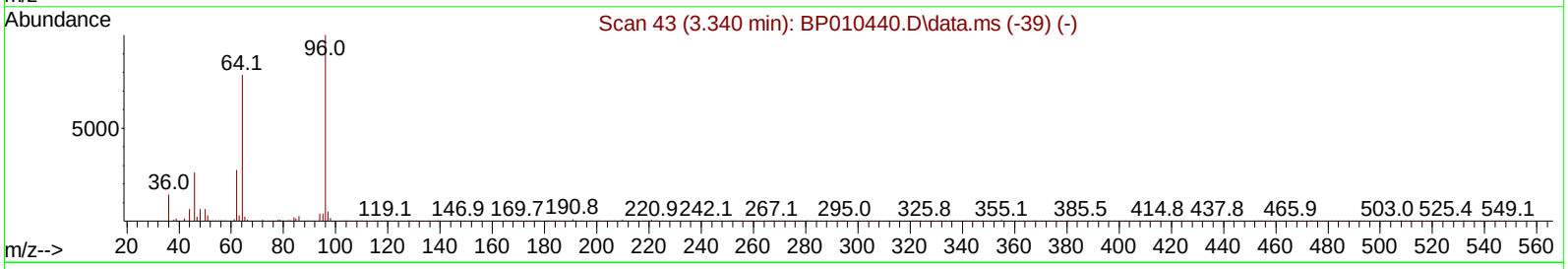

TIC: BP010450.D\data.ms

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

3.340min (-0.006) 4.76 ng/uL m

response 15350

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.67#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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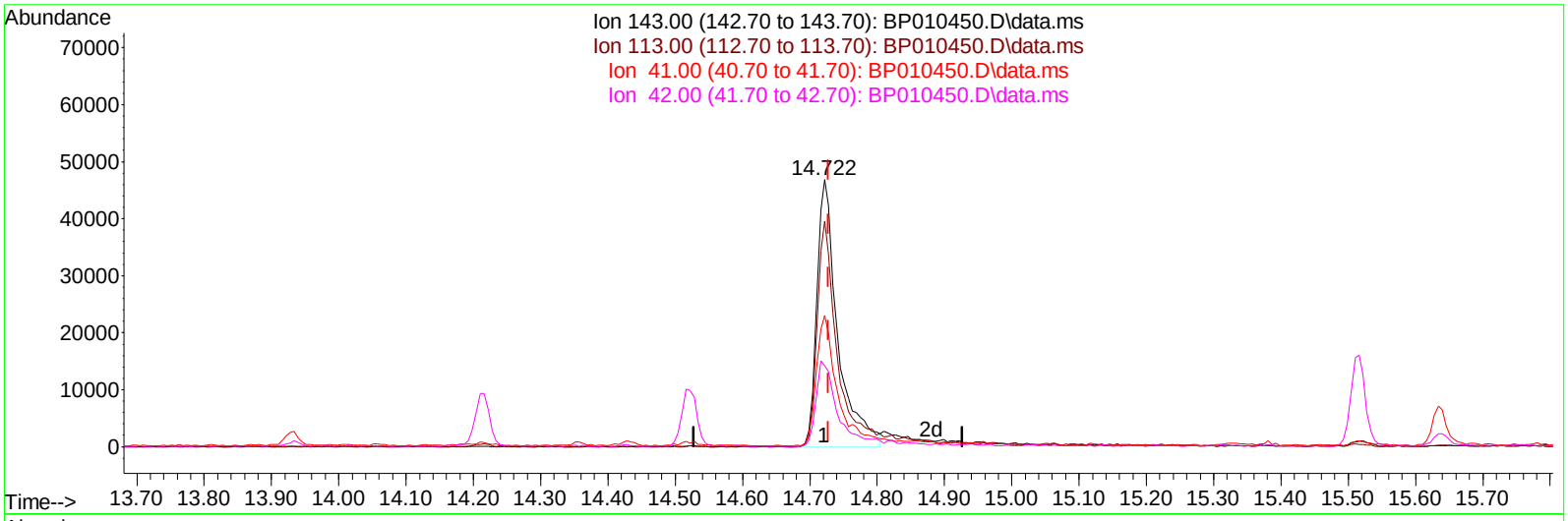

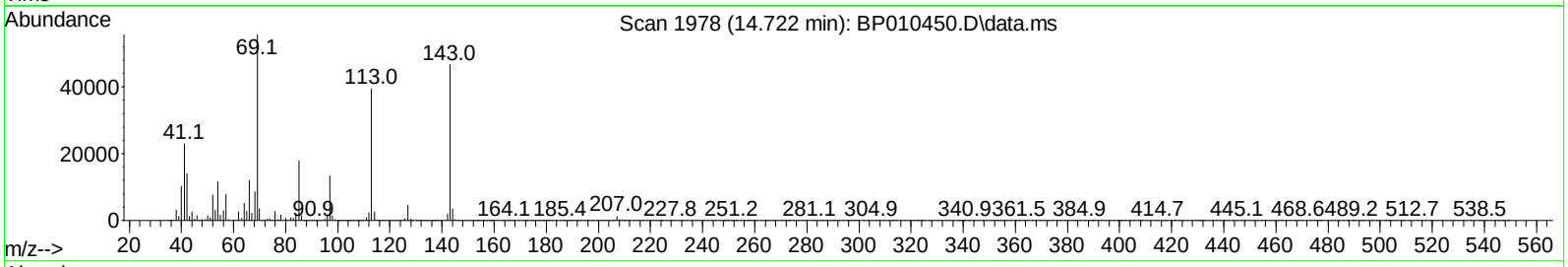

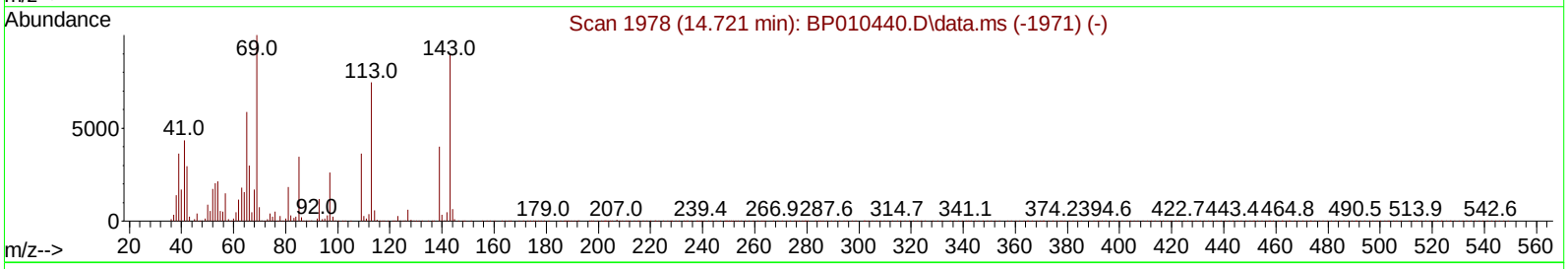

TIC: BP010450.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 18.32 ng/ul

response 100374

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.48#
41.00	23.20	49.20#
42.00	18.00	30.35#

Quantitation Report (Qedit)

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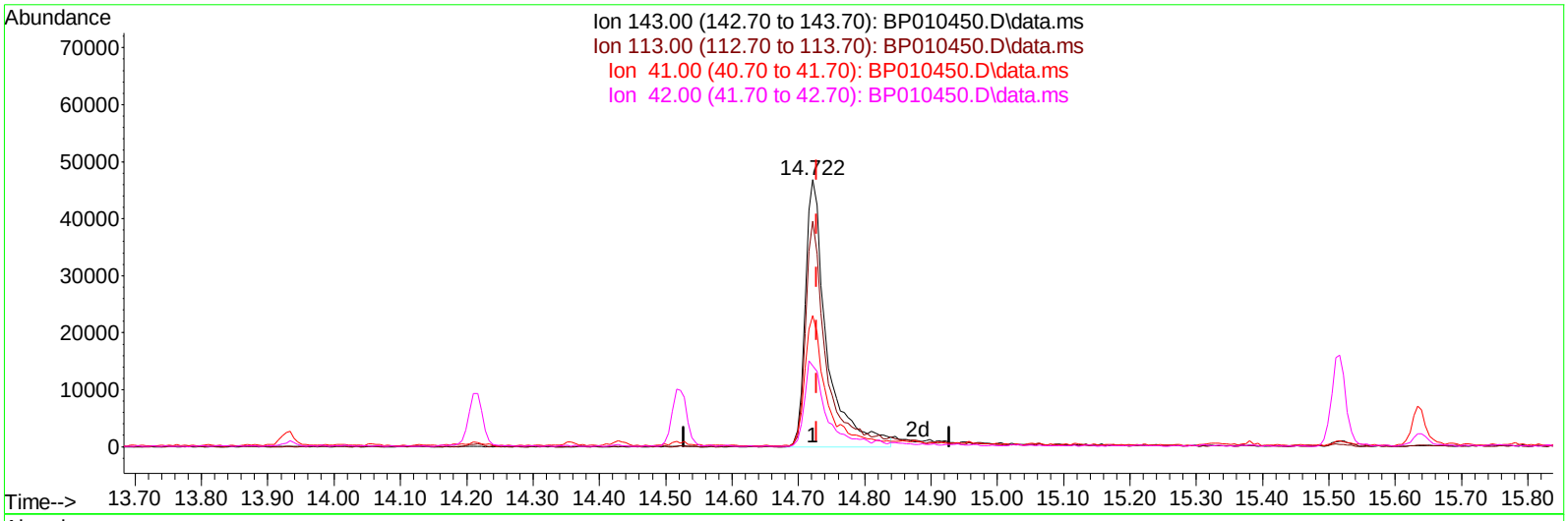

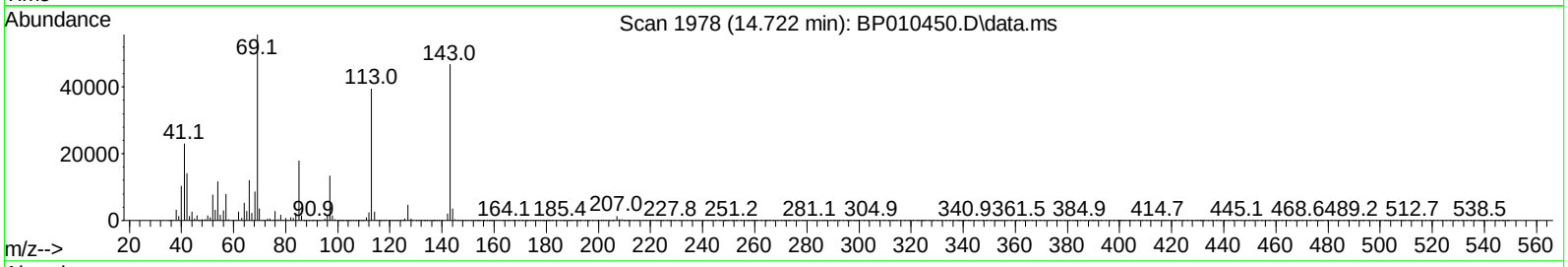

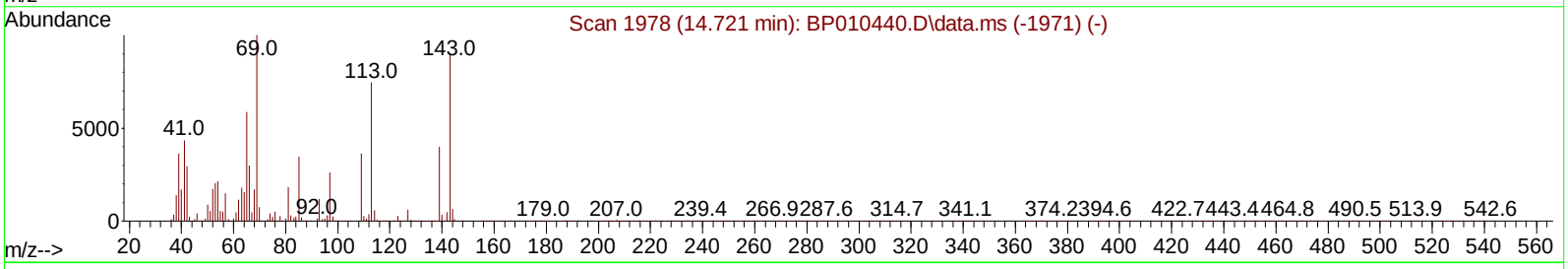

TIC: BP010450.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 19.14 ng/ul m

response 104886

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	84.48#
41.00	23.20	49.20#
42.00	18.00	30.35#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	137377	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	616720	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	417798	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	913946	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	860301	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	735050	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	15350m	4.763	ng/uL	0.00
4) Pyridine-d5	3.764	84	104629	11.433	ng/ul	0.00
7) Phenol-d5	7.058	99	291209	25.256	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	211171	27.259	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	240784	27.291	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	236821	24.168	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	131214	27.491	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.775	143	141996	28.072	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	241863	25.288	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	342321	24.071	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	894220	29.050	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1015357	28.597	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	104886m	19.143	ng/ul	0.00
60) Fluorene-d10	15.516	176	769886	28.749	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	100887	18.875	ng/ul	-0.01
73) Anthracene-d10	17.375	188	1222959	29.551	ng/ul	0.00
81) Pyrene-d10	19.604	212	1504687	32.934	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1154808	31.287	ng/ul	-0.01

Target Compounds Qvalue

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

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