Quantitation Report (Qedit)

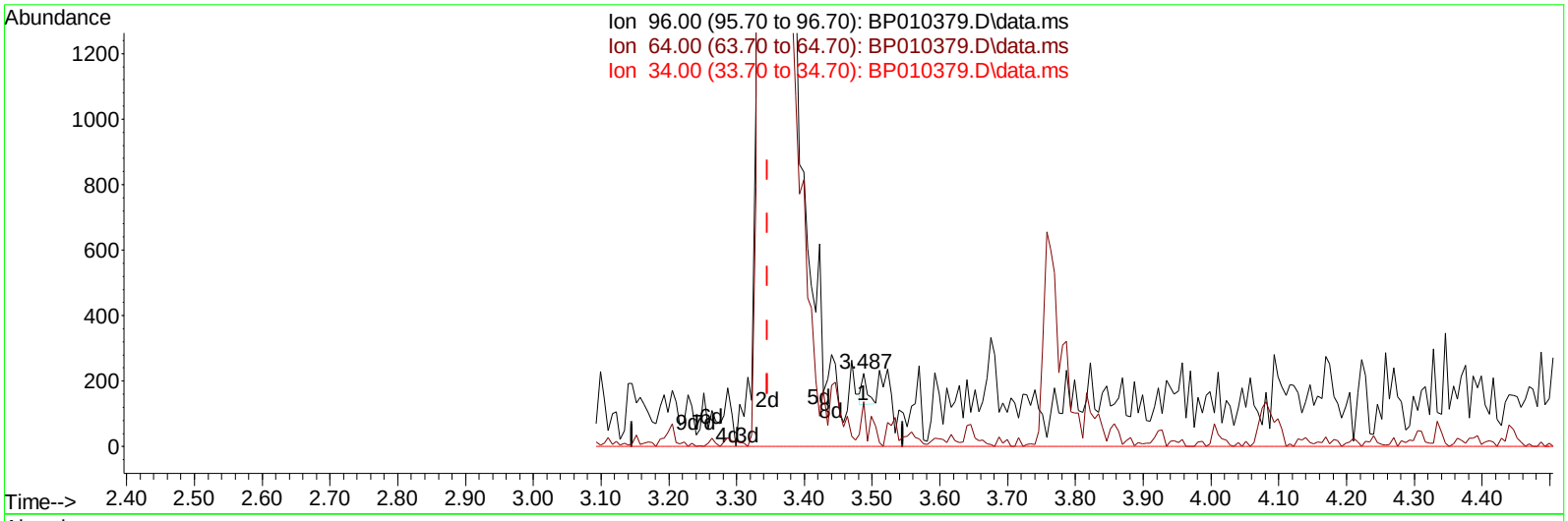
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010379.D
Acq On : 17 May 2022 04:15
Operator : CG/JU
Sample : PB144810BL
Misc :
ALS Vial : 27 Sample Multiplier: 1

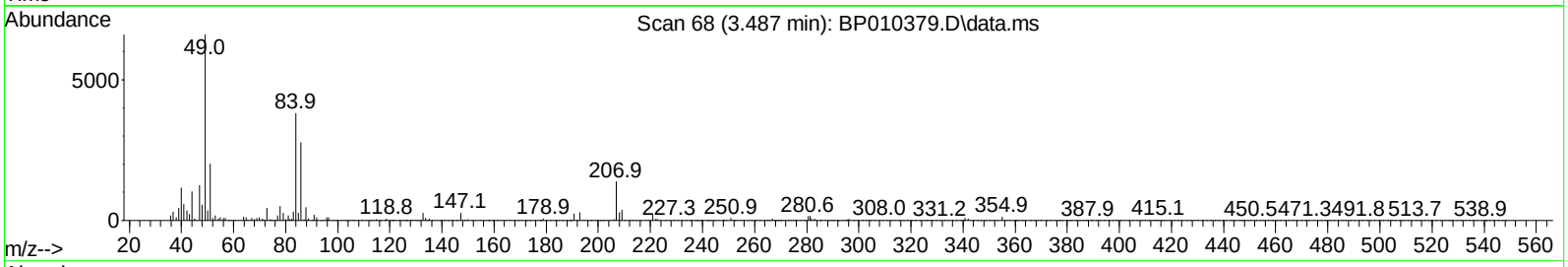
Quant Time: May 17 05:03:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 13 15:04:31 2022
Response via : Initial Calibration

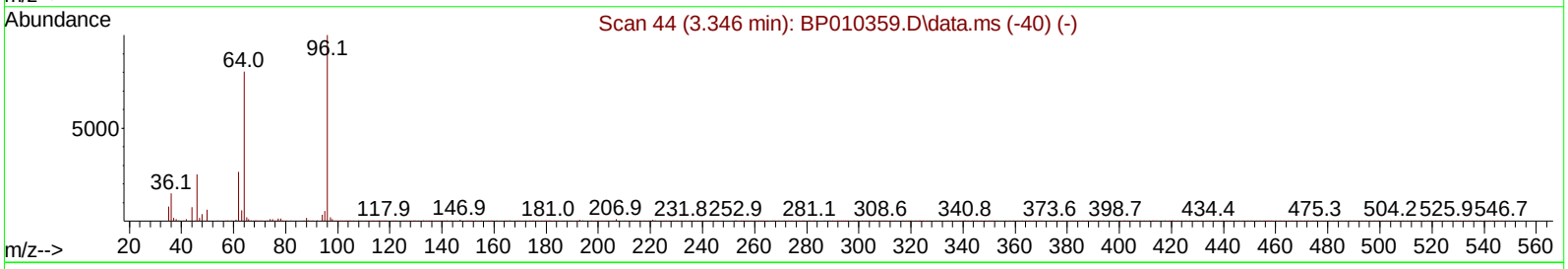
Instrument :
BNA_P
ClientSampleId :
SBLK810

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/17/2022
Supervised By :mohammad ahmed 05/18/2022







TIC: BP010379.D\data.ms

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

3.487min (+ 0.141) 0.01 ng/uL

response 52

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	61.43
34.00	0.00	0.00
0.00	0.00	0.00

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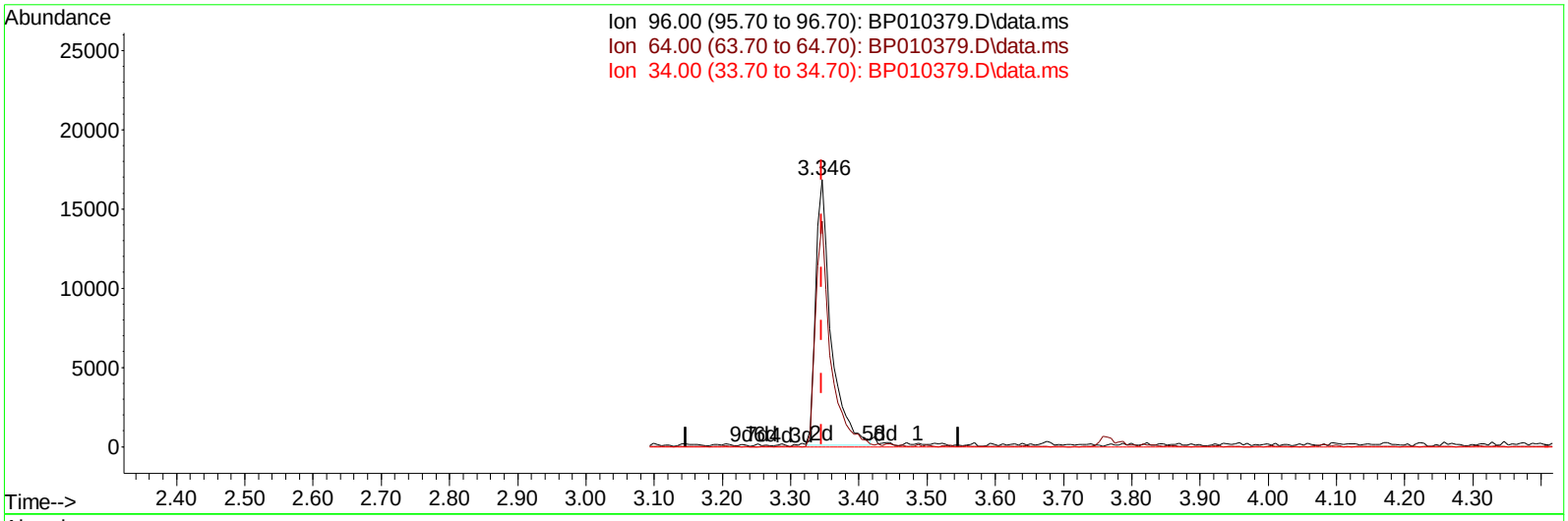

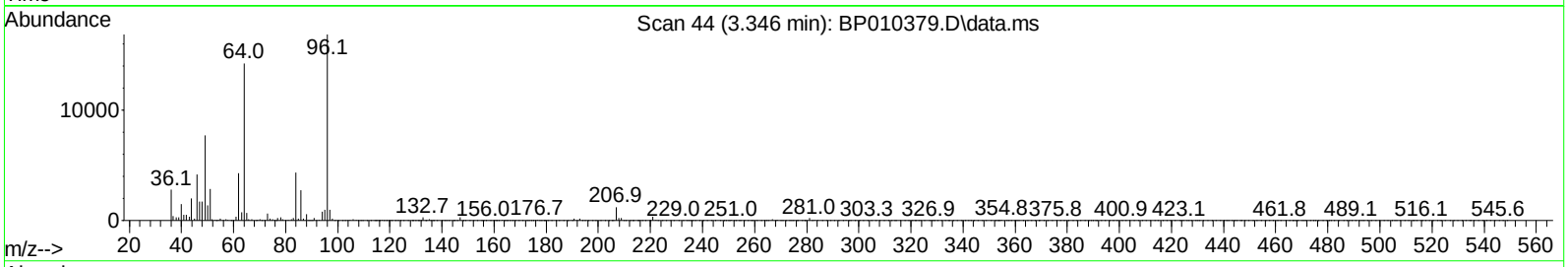

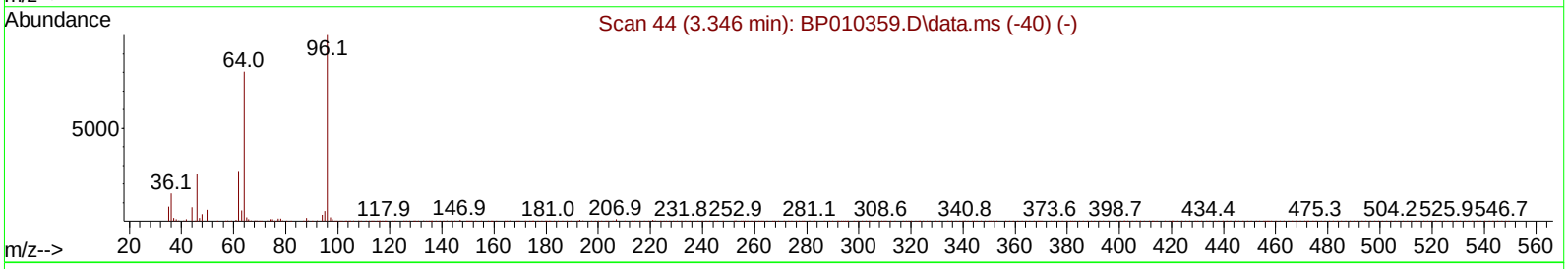

TIC: BP010379.D\data.ms

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

3.346min (-0.000) 7.16 ng/uL m

response 26367

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	84.41#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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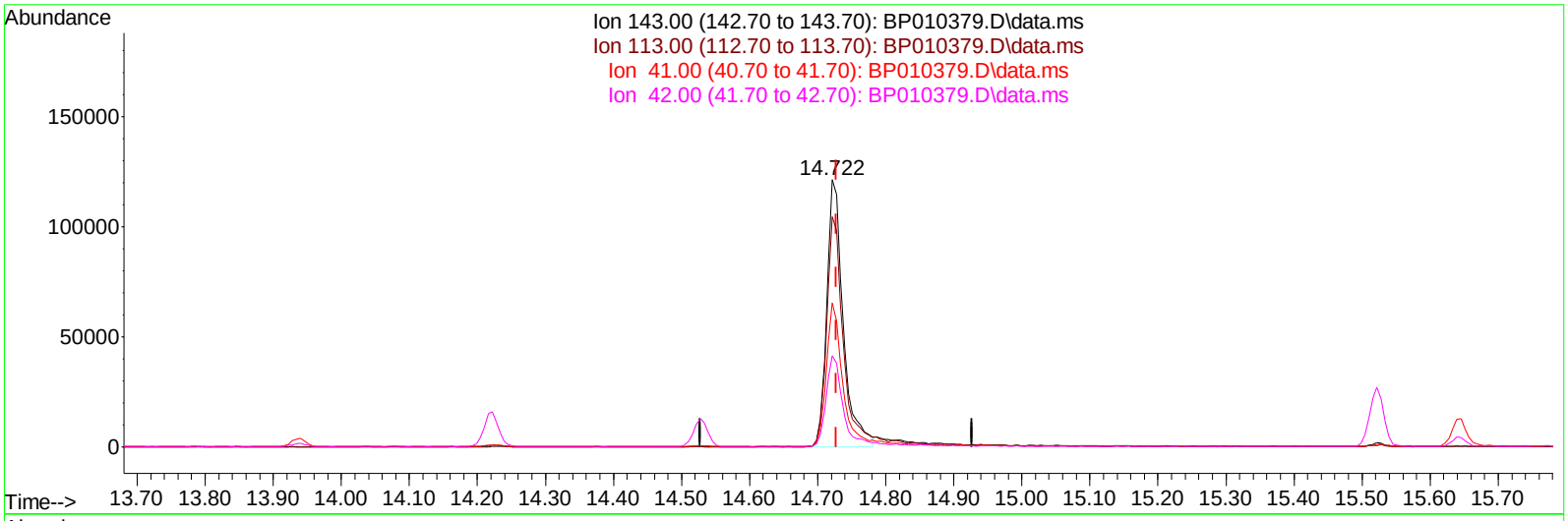

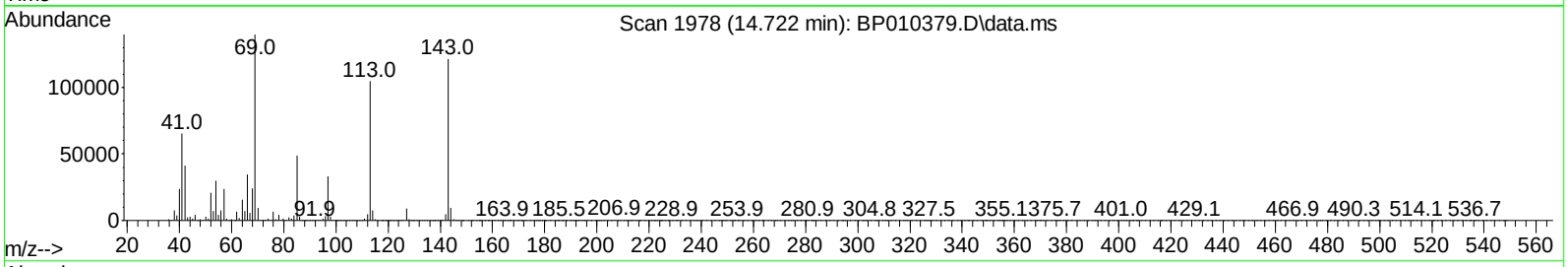

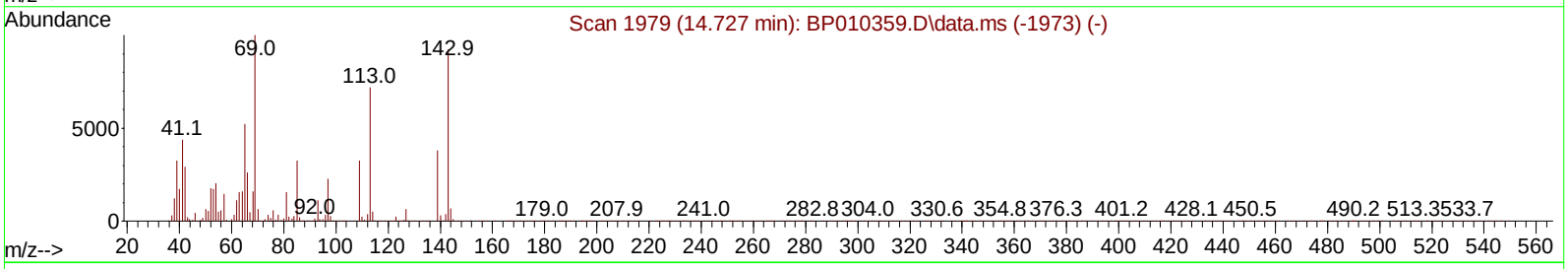

TIC: BP010379.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 31.73 ng/ul

response 202338

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.15#
41.00	23.20	53.78#
42.00	18.00	34.18#

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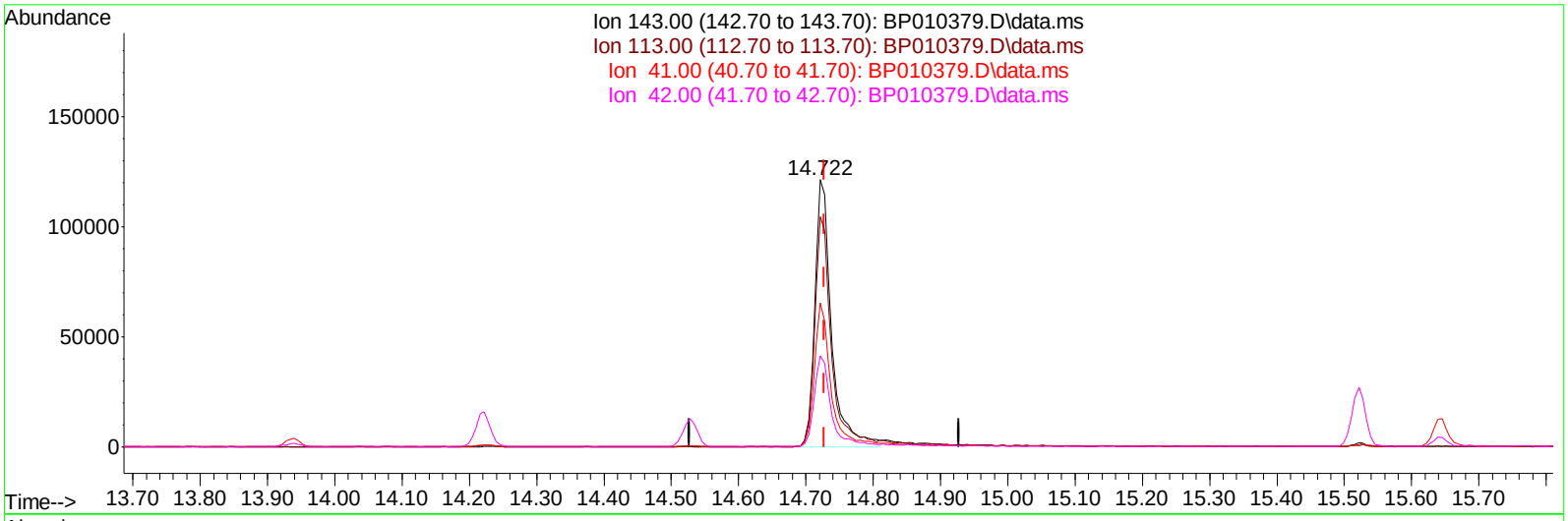

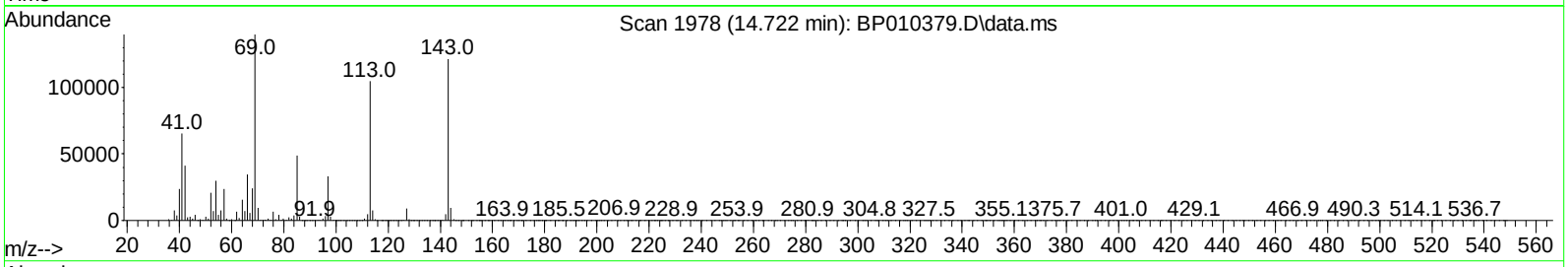

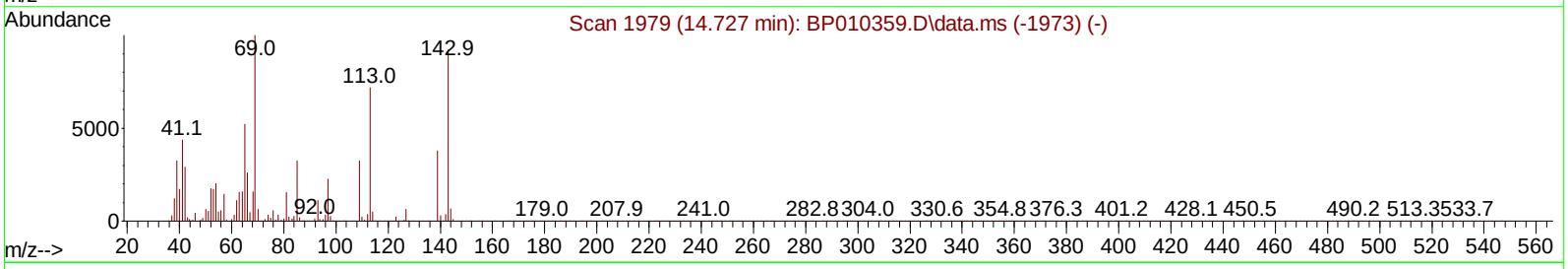

TIC: BP010379.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 32.72 ng/ul m

response 208667

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.15#
41.00	23.20	53.78#
42.00	18.00	34.18#

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	7.899	152	156965	20.000	ng/ul	0.00
20) Naphthalene-d8	10.698	136	698477	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	486288	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1058417	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	1004312	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	830898	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	26367m	7.160	ng/uL	0.00
4) Pyridine-d5	3.758	84	381981	36.532	ng/ul	0.00
7) Phenol-d5	7.063	99	489925	37.188	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	346061	39.096	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	397045	39.386	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	428296	38.254	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	212590	39.327	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	235902	41.178	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	404858	37.375	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	593756	36.865	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	1450871	40.495	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1629045	39.419	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	208667m	32.721	ng/ul	0.00
60) Fluorene-d10	15.522	176	1235165	39.627	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	190505	30.776	ng/ul	0.00
73) Anthracene-d10	17.380	188	1954332	40.777	ng/ul	0.00
81) Pyrene-d10	19.616	212	2350741	44.074	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1793253	42.981	ng/ul	0.00
Target Compounds						Qvalue

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

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