Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
Data File : BP010364.D
Acq On : 16 May 2022 19:12
Operator : CG/JU
Sample : N2831-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: May 16 23:22:05 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 :
YB2K1

Manual IntegrationsAPPROVED

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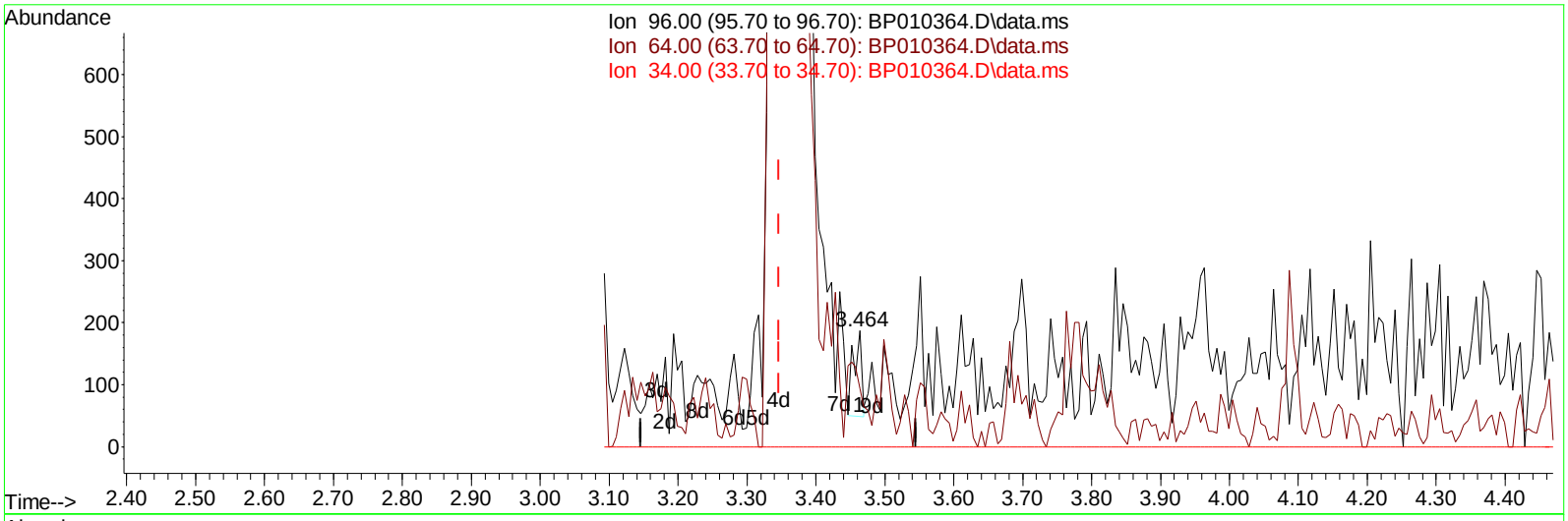
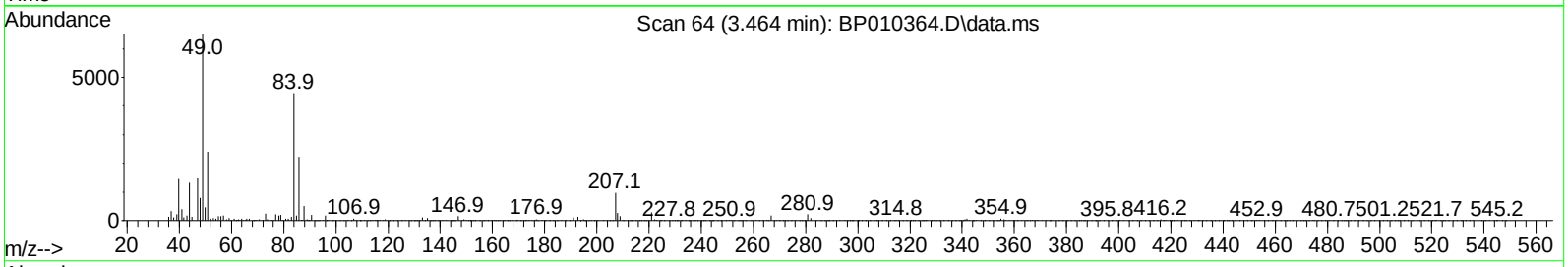
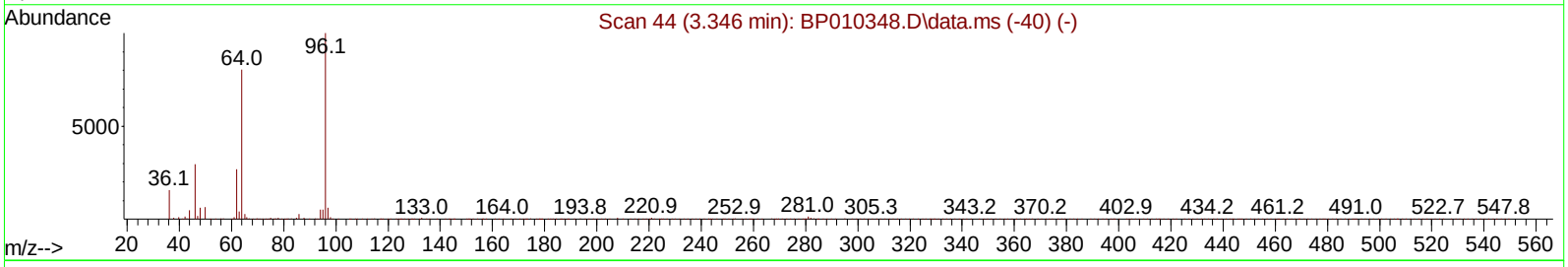
Supervised By :mohammad ahmed 05/18/2022







TIC: BP010364.D\data.ms

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

3.464min (+ 0.118) 0.04 ng/uL

response 118

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	50.27
34.00	0.00	0.00
0.00	0.00	0.00

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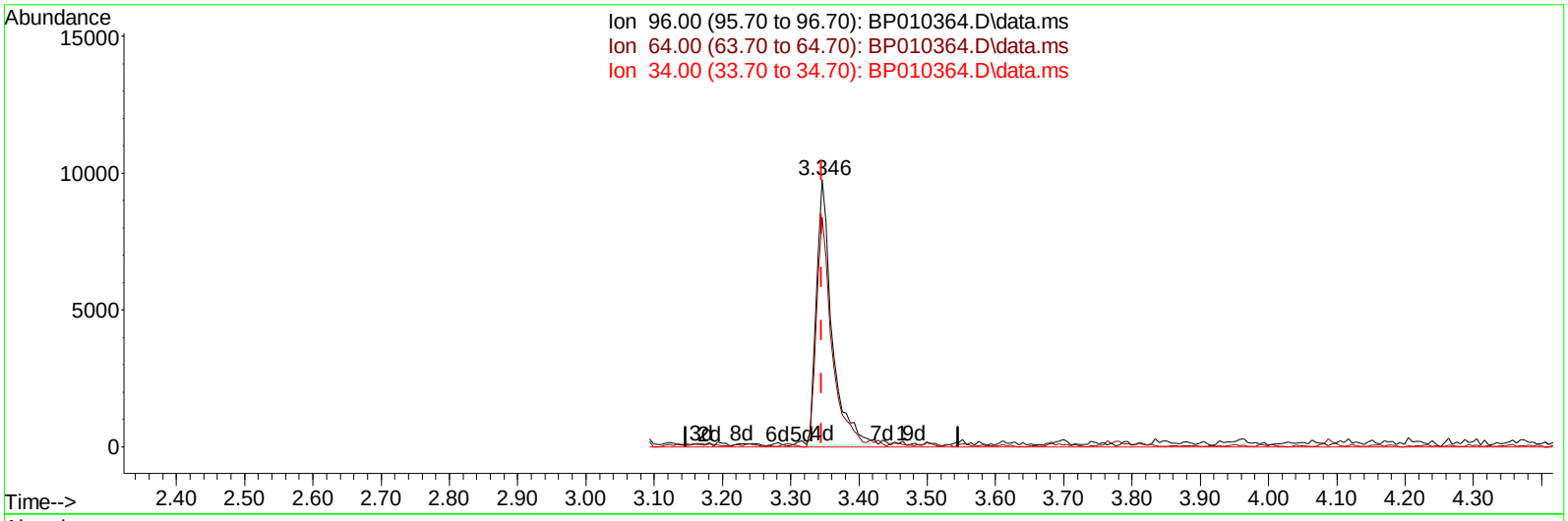

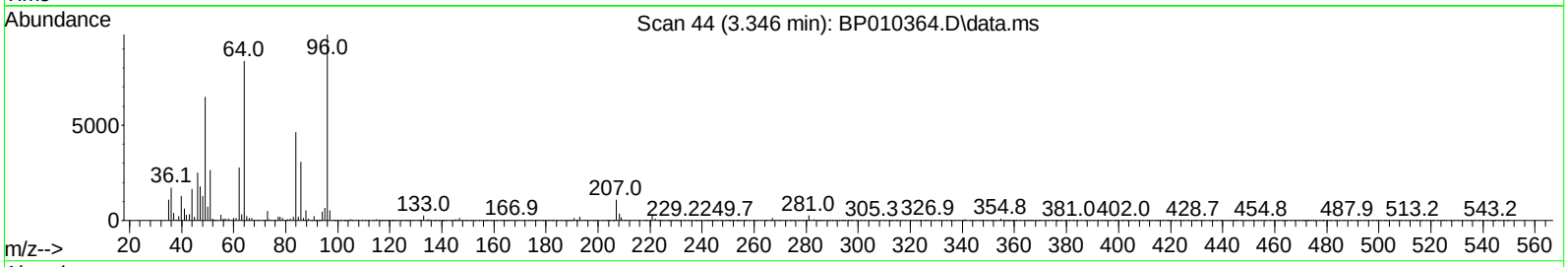

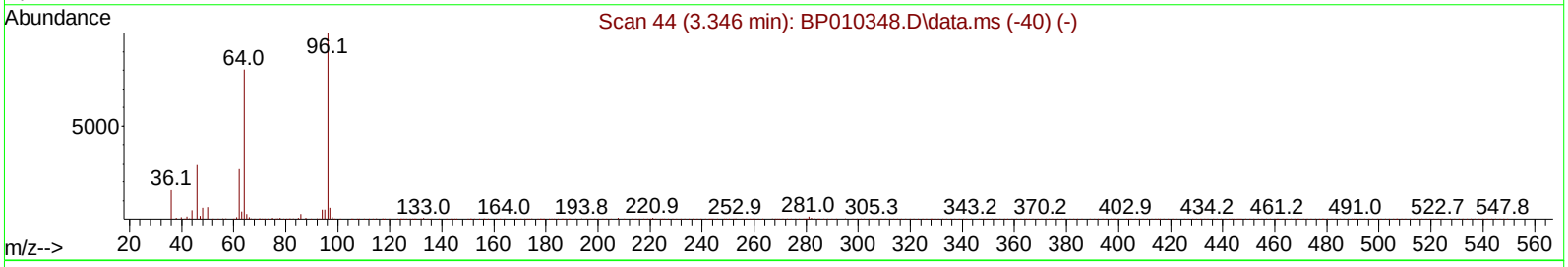

TIC: BP010364.D\data.ms

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

3.346min (+ 0.000) 4.62 ng/uL m

response 15514

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	85.84#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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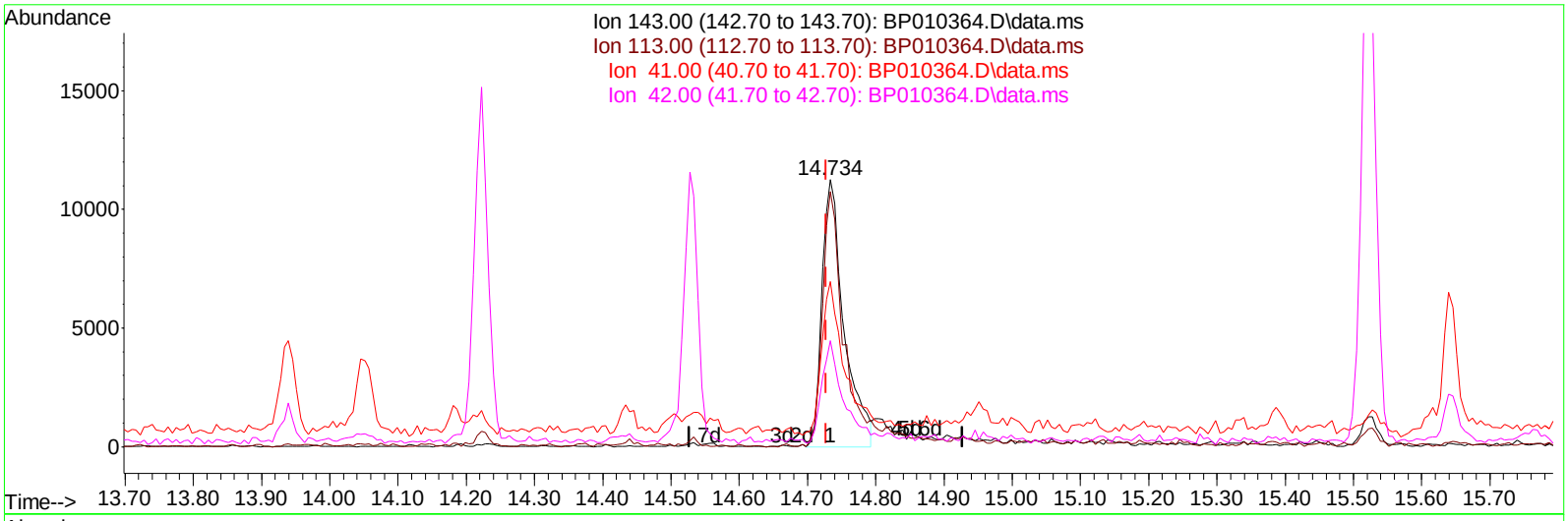

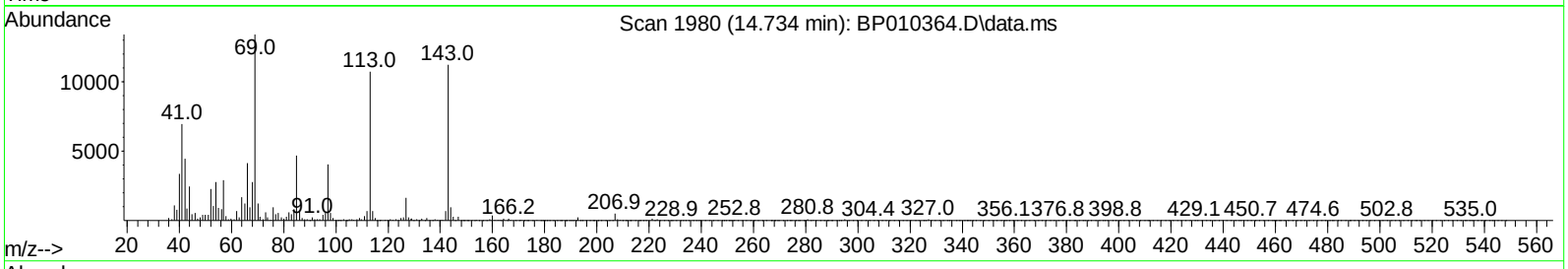

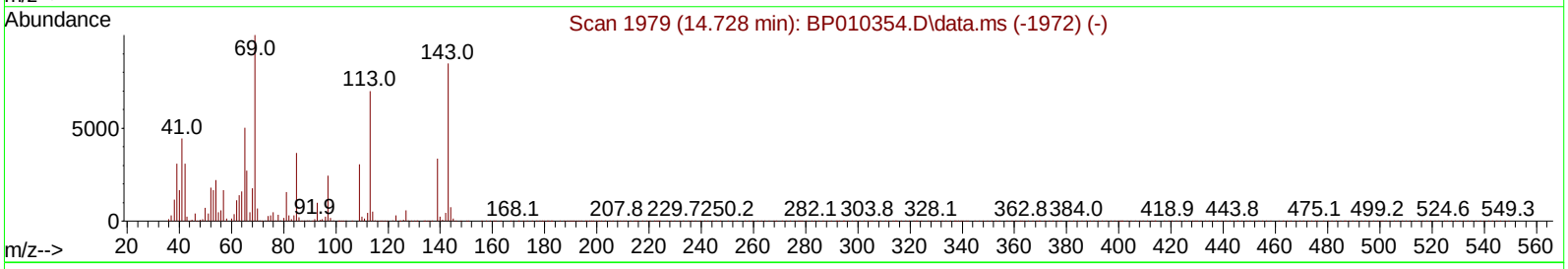

TIC: BP010364.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.66 ng/ul

response 25331

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	95.55#
41.00	23.20	61.90#
42.00	18.00	39.72#

Quantitation Report (Qedit)

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Instrument :
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ClientSampleId :
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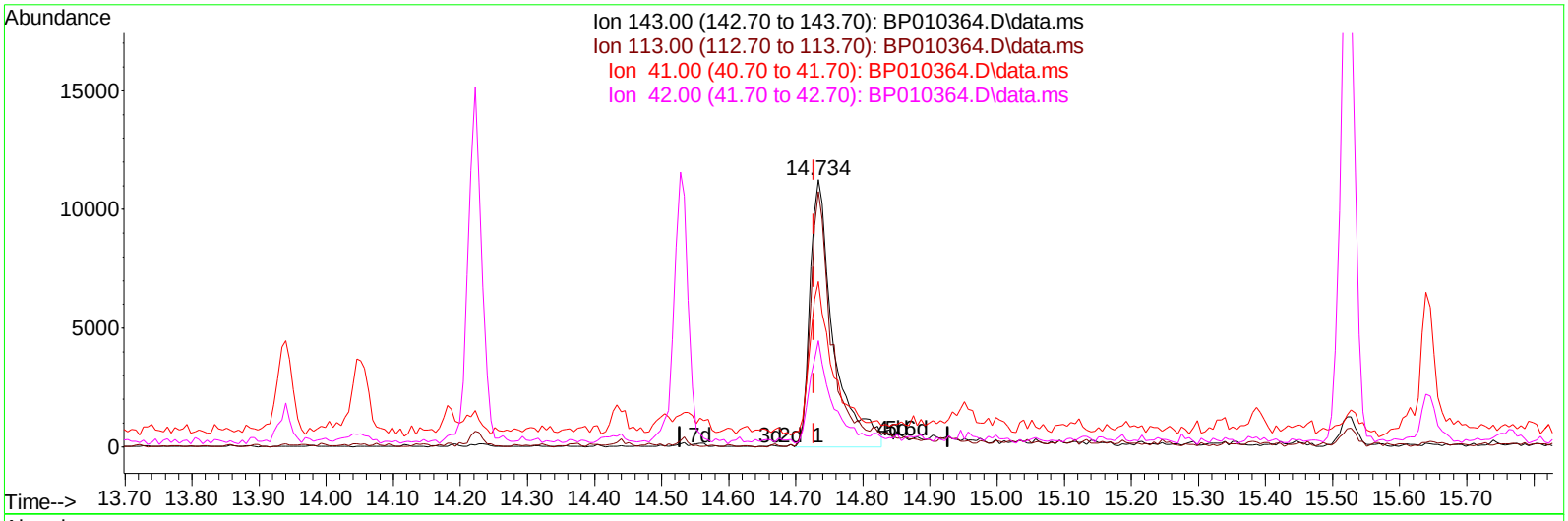

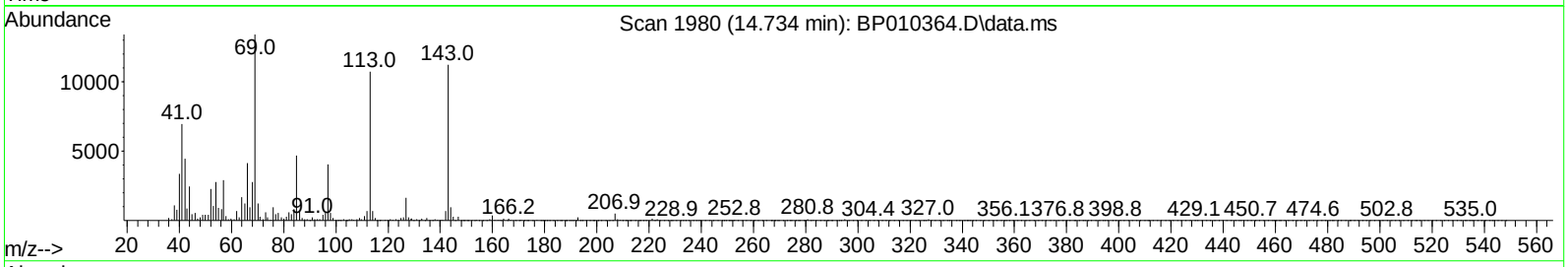

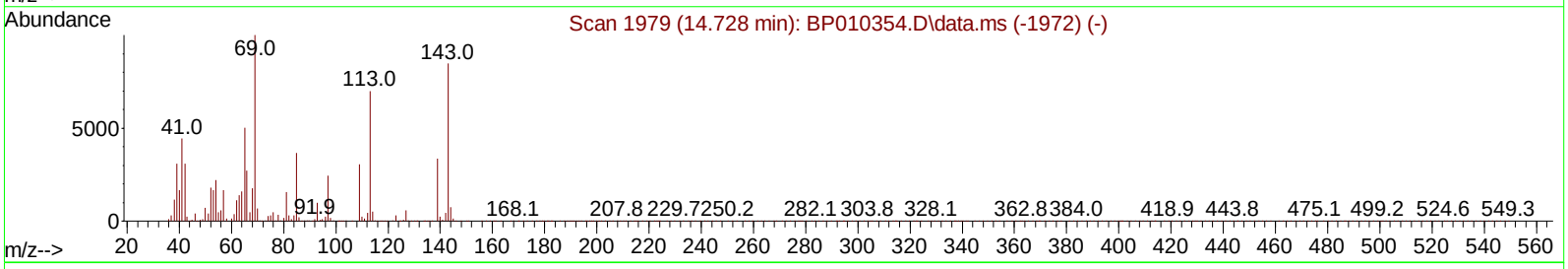

TIC: BP010364.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 5.03 ng/ul m

response 27325

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	95.55#
41.00	23.20	61.90#
42.00	18.00	39.72#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	143277	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	615259	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	414561	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	916723	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	925311	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	819421	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	15514m	4.615	ng/uL	0.00
4) Pyridine-d5	3.770	84	103808	10.877	ng/ul	0.00
7) Phenol-d5	7.064	99	86909	7.227	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	275410	34.087	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	248351	26.989	ng/ul	0.00
15) 4-Methylphenol-d8	8.605	113	181063	17.717	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	168904	35.471	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	160507	31.807	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	323504	33.904	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	447044	31.510	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1211857	39.676	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1379194	39.147	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	27325m	5.026	ng/ul	0.00
60) Fluorene-d10	15.522	176	1084351	40.808	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	77186	14.397	ng/ul	0.00
73) Anthracene-d10	17.381	188	1745035	42.038	ng/ul	0.00
81) Pyrene-d10	19.616	212	2159669	43.949	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1749959	42.530	ng/ul	0.00

Target Compounds Qvalue

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

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