Quantitation Report (Qedit)

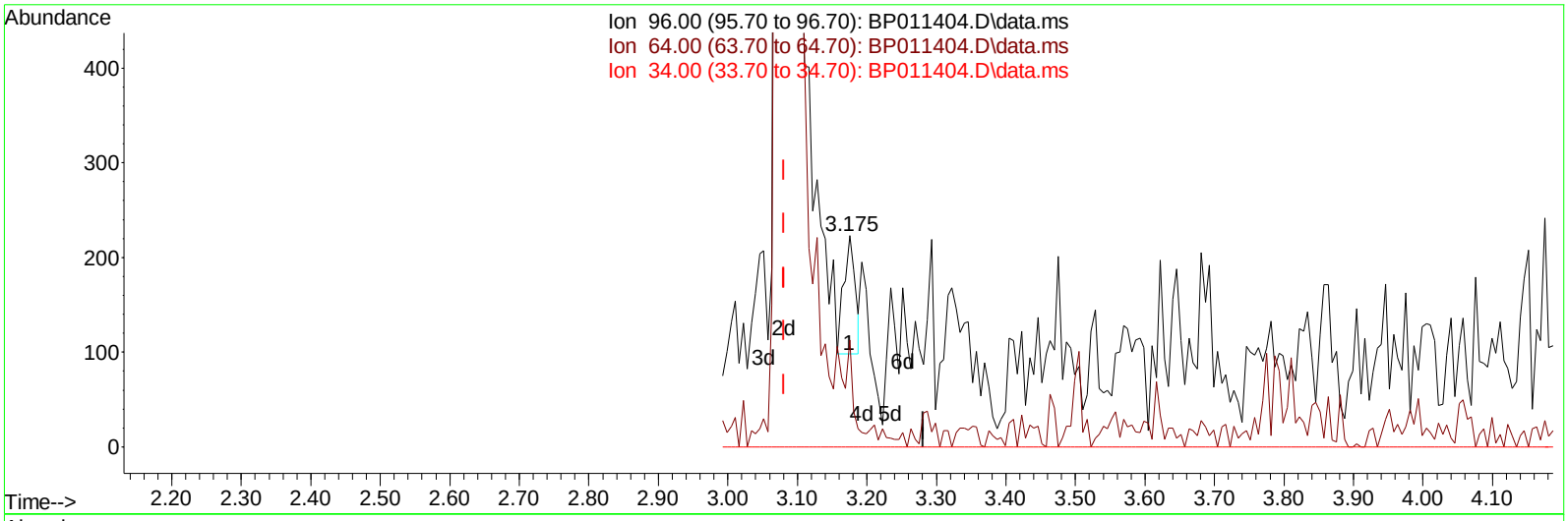
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081122\
Data File : BP011404.D
Acq On : 12 Aug 2022 01:25
Operator : CG/JU
Sample : N4128-09
Misc :
ALS Vial : 27 Sample Multiplier: 1

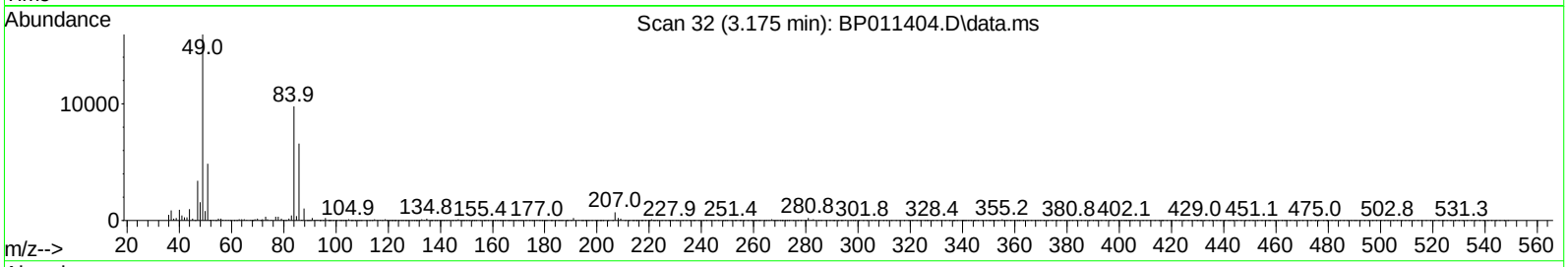
Quant Time: Aug 12 02:25:24 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 11 23:05:53 2022
Response via : Initial Calibration

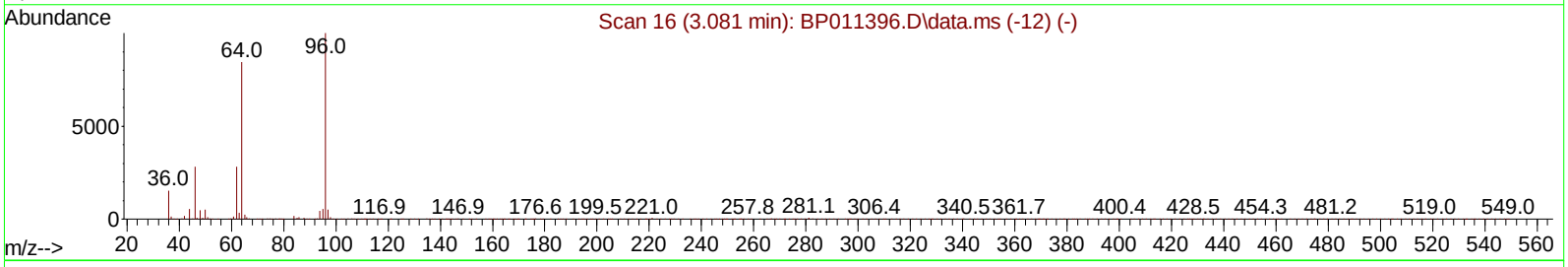
Instrument :
BNA_P
ClientSampleId :
BGBT6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/12/2022
Supervised By :Sohil Jodhani 08/13/2022







TIC: BP011404.D\data.ms

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

3.175min (+ 0.094) 0.05 ng/uL

response 142

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	50.67
34.00	0.00	0.00
0.00	0.00	0.00

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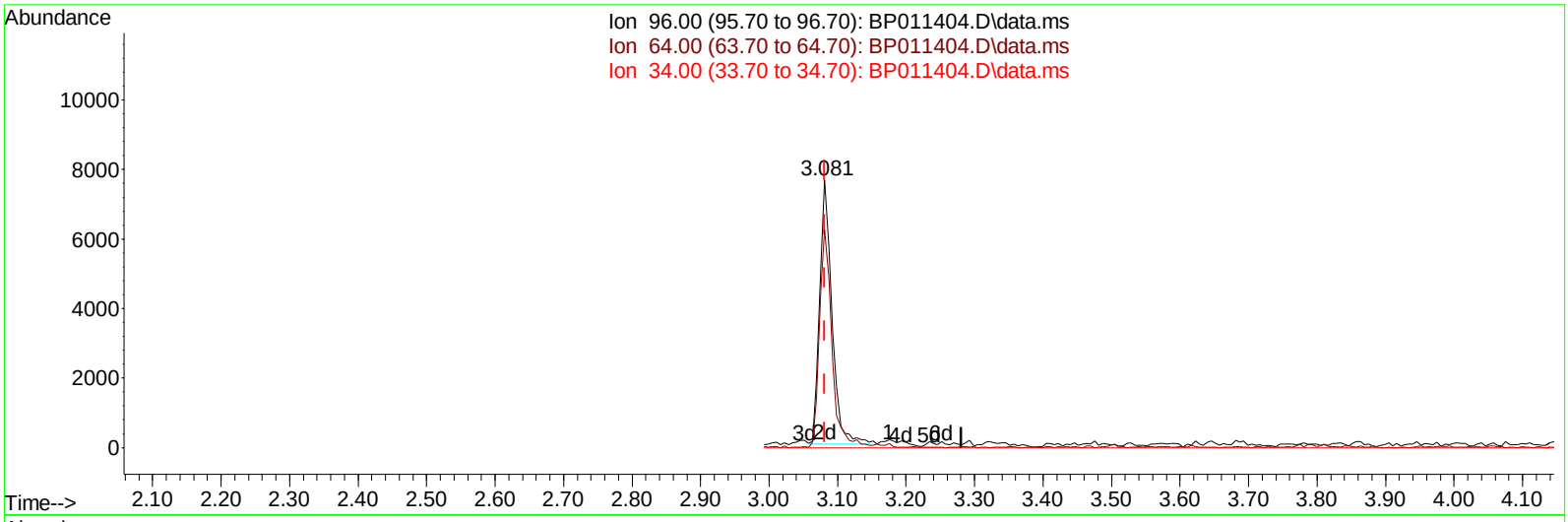

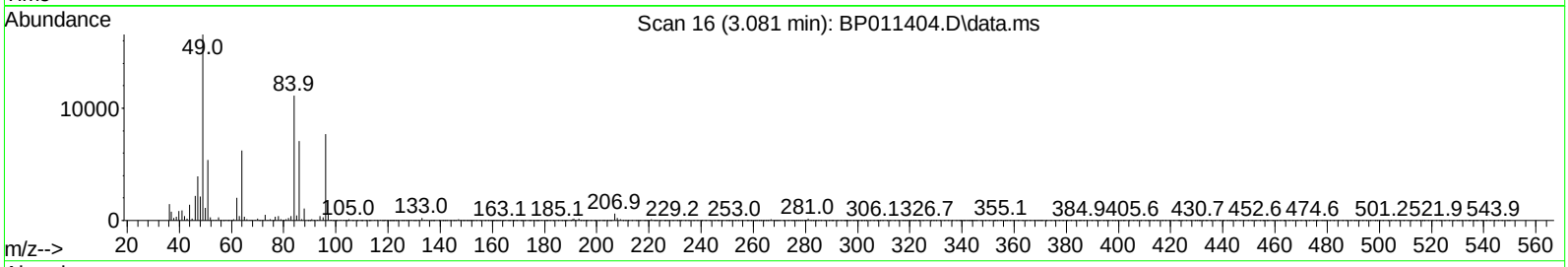

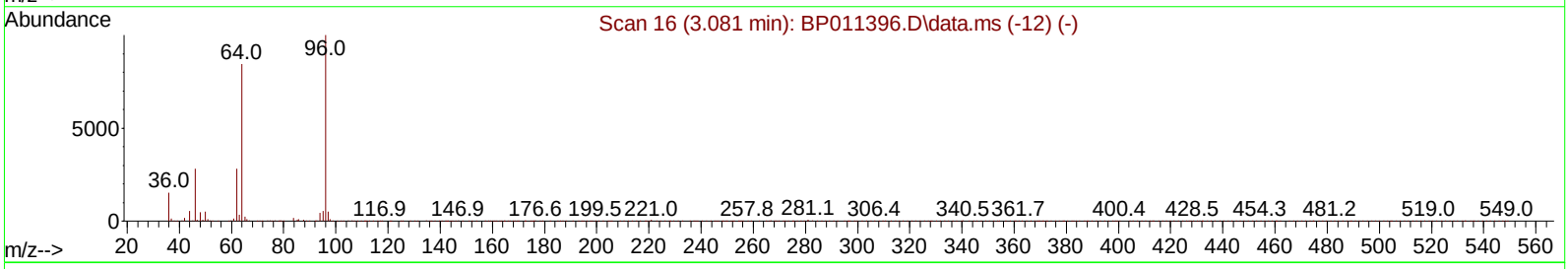

TIC: BP011404.D\data.ms

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

3.081min (-0.000) 3.27 ng/uL m

response 9455

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	81.36#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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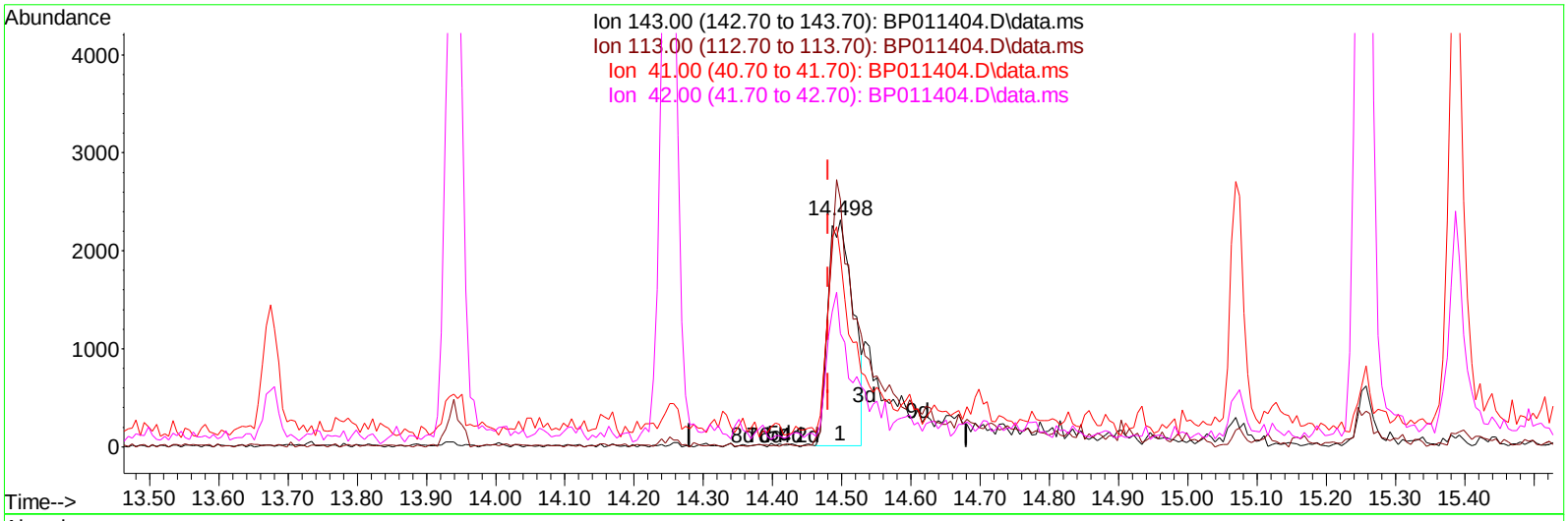

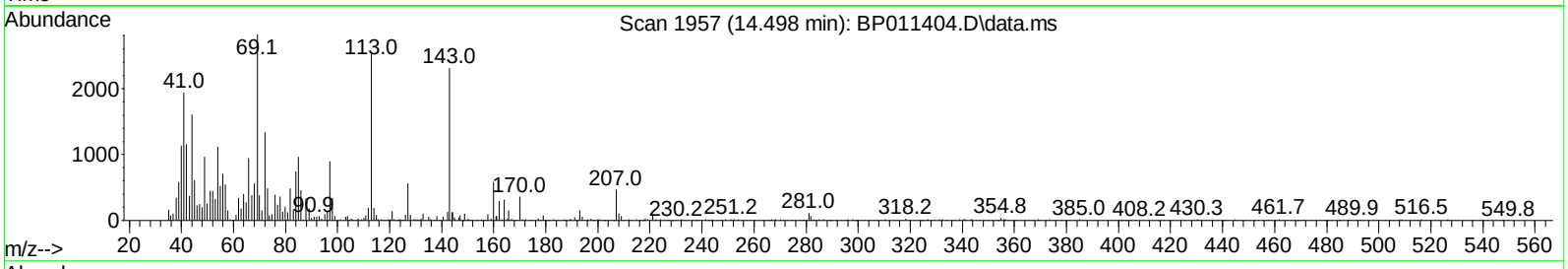

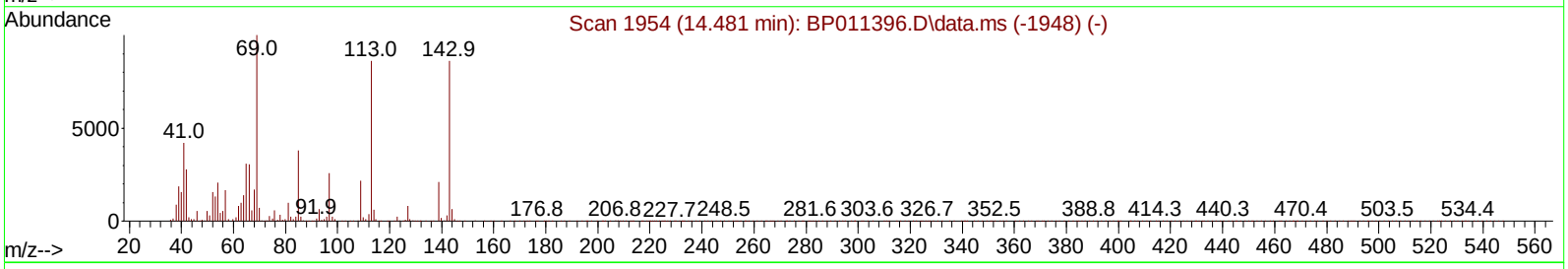

TIC: BP011404.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 2.19 ng/ul

response 5774

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	108.93#
41.00	54.40	83.82#
42.00	35.50	49.89#

Quantitation Report (Qedit)

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Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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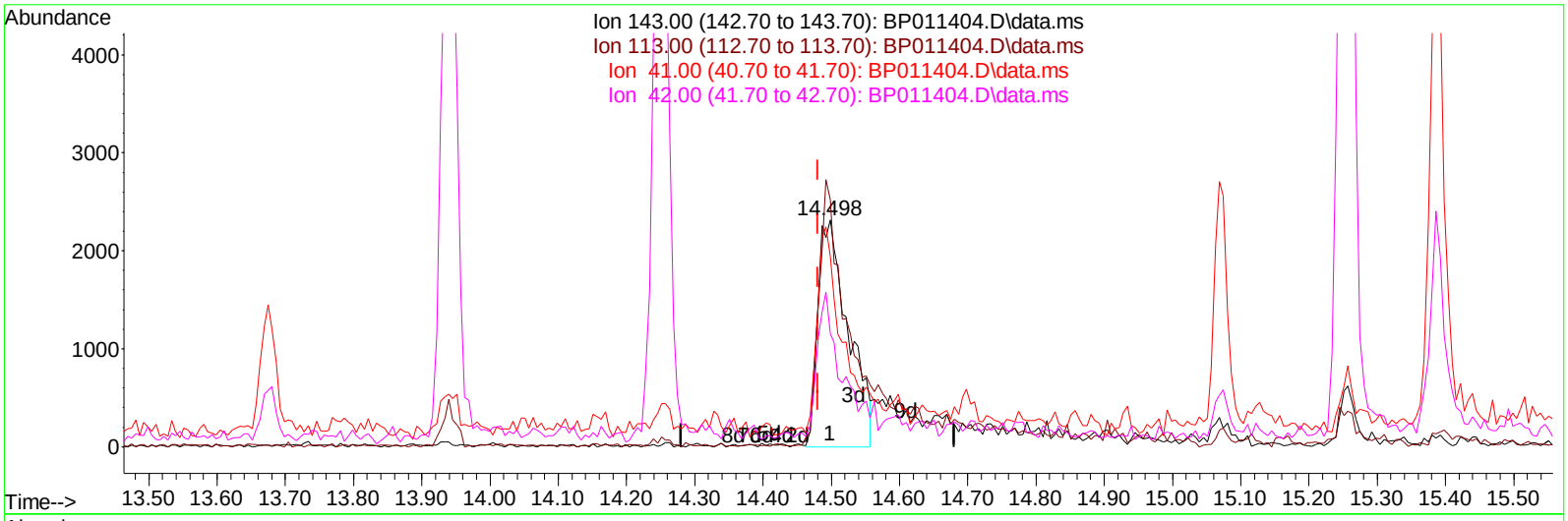

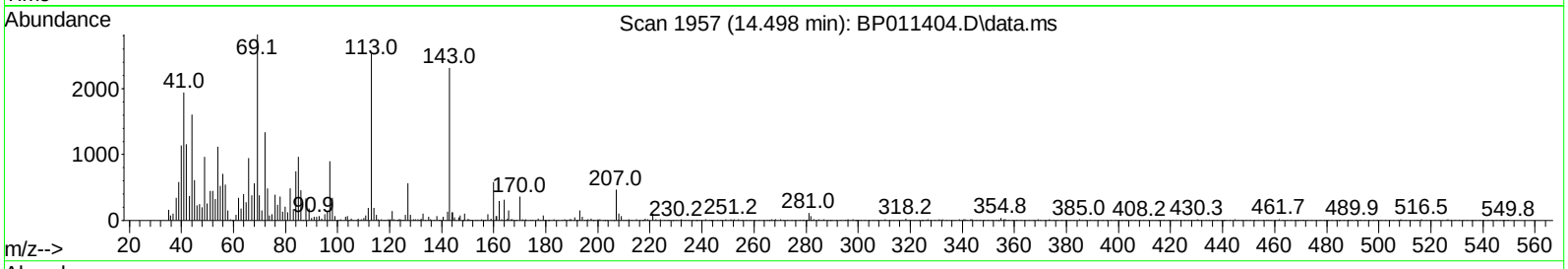

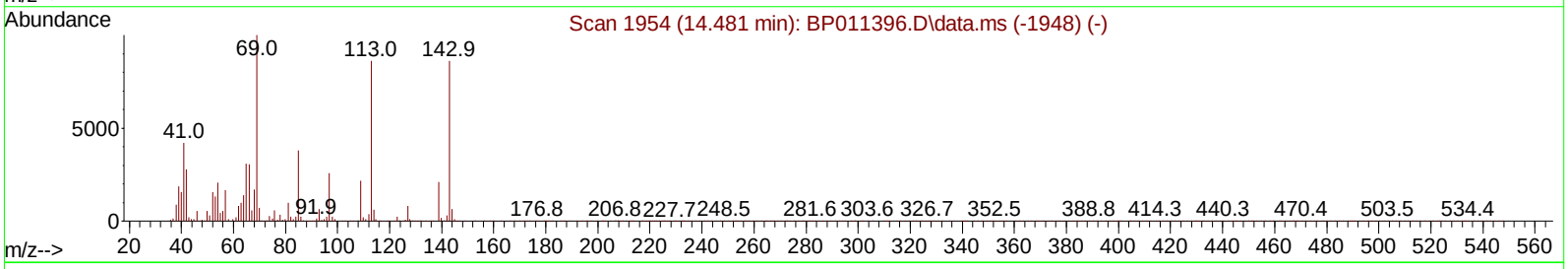

TIC: BP011404.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.018) 2.74 ng/ul m

response 7209

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	108.93#
41.00	54.40	83.82#
42.00	35.50	49.89#

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

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Supervised By :Sohil Jodhani 08/13/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.575	152	88701	20.000	ng/ul		0.00
20) Naphthalene-d8	10.369	136	371649	20.000	ng/ul		0.00
38) Acenaphthene-d10	14.251	164	207921	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.022	188	436607	20.000	ng/ul		0.00
79) Chrysene-d12	21.157	240	426125	20.000	ng/ul	#	0.00
88) Perylene-d12	23.462	264	457577	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.081	96	9455m	3.268	ng/uL		0.00
4) Pyridine-d5	3.499	84	37234	5.101	ng/ul		0.01
7) Phenol-d5	6.763	99	41321	5.264	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	6.922	67	140280	25.204	ng/ul		0.00
11) 2-Chlorophenol-d4	7.110	132	127088	20.634	ng/ul		0.00
15) 4-Methylphenol-d8	8.299	113	76649	12.436	ng/ul		0.00
21) Nitrobenzene-d5	8.740	128	76651	30.305	ng/ul		0.00
24) 2-Nitrophenol-d4	9.457	143	72701	31.394	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	9.993	165	115364	24.669	ng/ul		0.00
31) 4-Chloroaniline-d4	10.522	131	147298	18.016	ng/ul		0.00
46) Dimethylphthalate-d6	13.675	166	453013	31.401	ng/ul		0.00
49) Acenaphthylene-d8	13.939	160	514679	29.637	ng/ul		0.00
54) 4-Nitrophenol-d4	14.498	143	7209m	2.736	ng/ul		0.02
60) Fluorene-d10	15.257	176	366139	29.677	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	53437	26.350	ng/ul		0.00
73) Anthracene-d10	17.122	188	656838	33.470	ng/ul		0.00
81) Pyrene-d10	19.386	212	775004	33.012	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.315	264	773094	33.502	ng/ul		0.00
Target Compounds							Qvalue

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

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