Quantitation Report (Qedit)

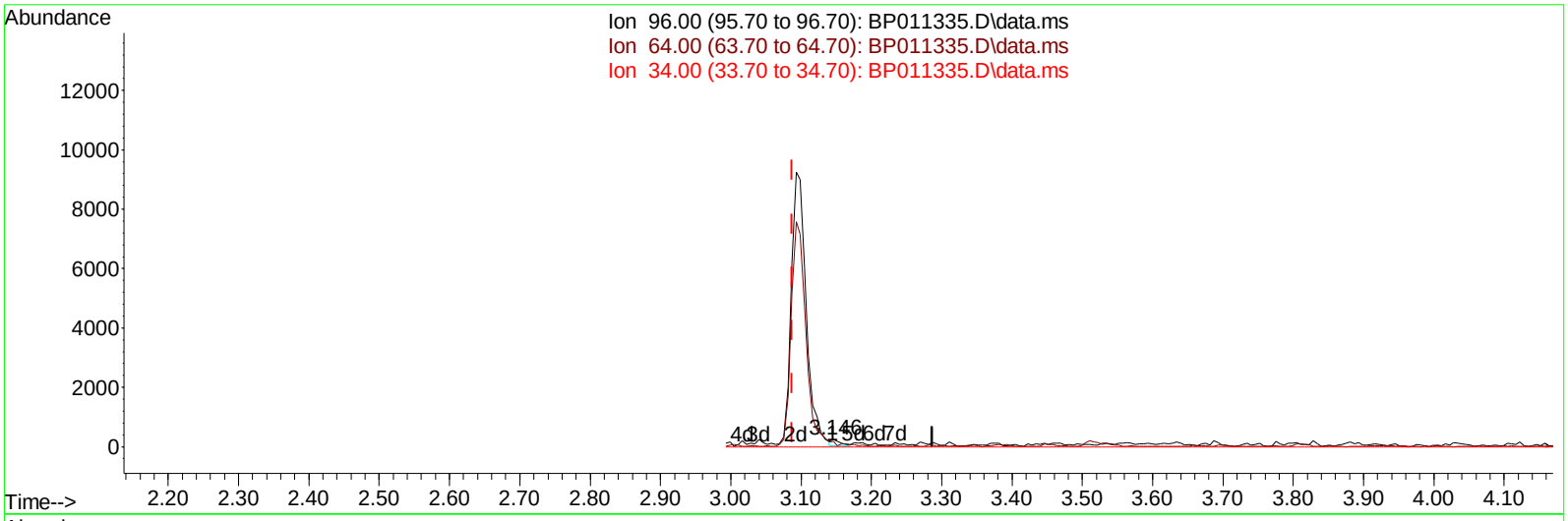
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011335.D
Acq On : 09 Aug 2022 18:24
Operator : CG/JU
Sample : N3900-13
Misc :
ALS Vial : 16 Sample Multiplier: 1

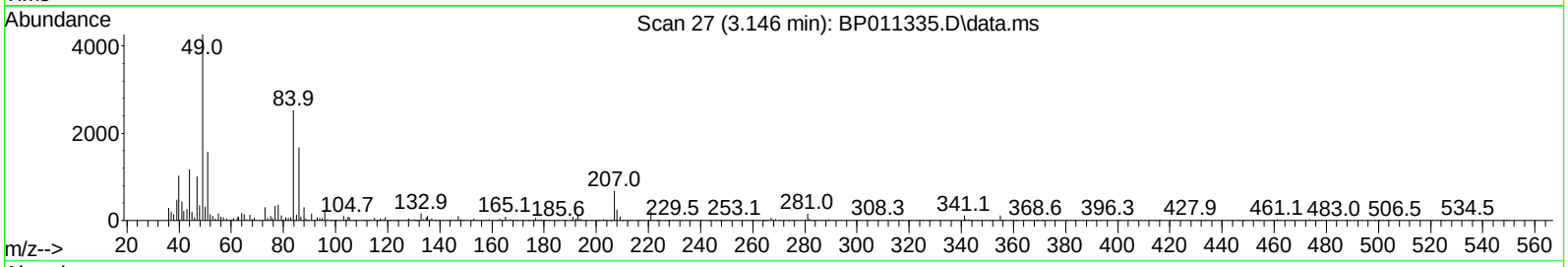
Quant Time: Aug 09 23:14:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

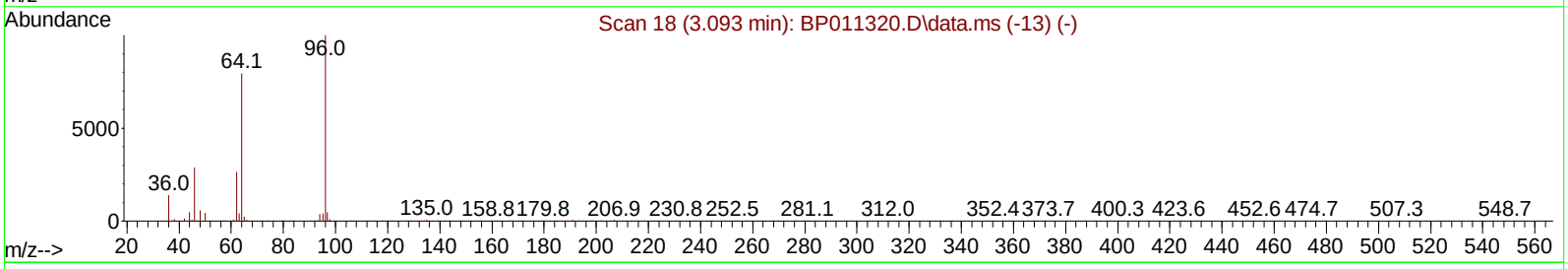
Instrument :
BNA_P
ClientSampleId :
C0C92

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011335.D\data.ms

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

3.146min (+ 0.059) 0.04 ng/uL

response 162

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	69.20
34.00	0.00	0.00
0.00	0.00	0.00

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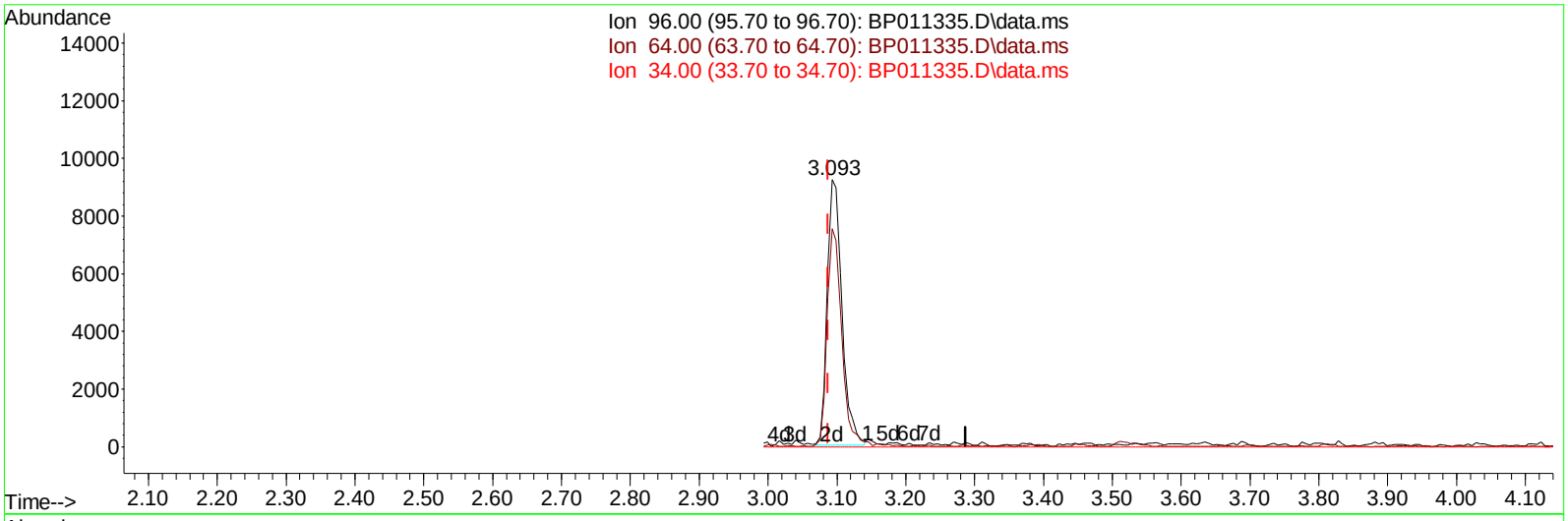

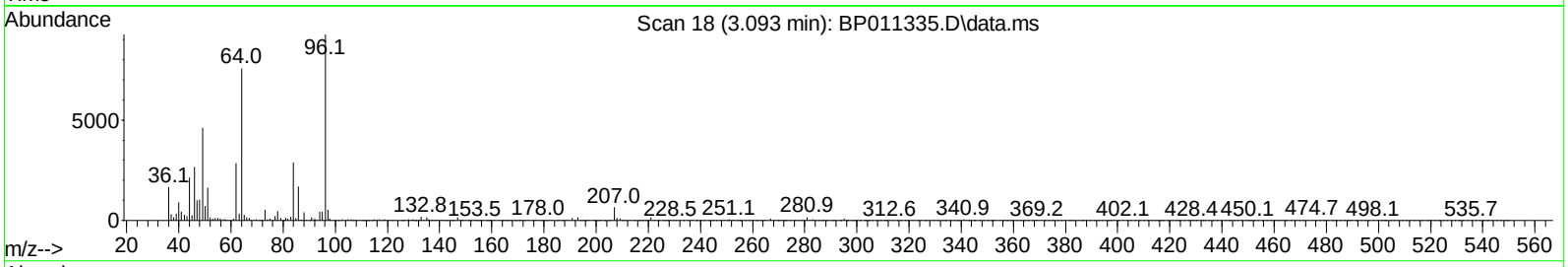

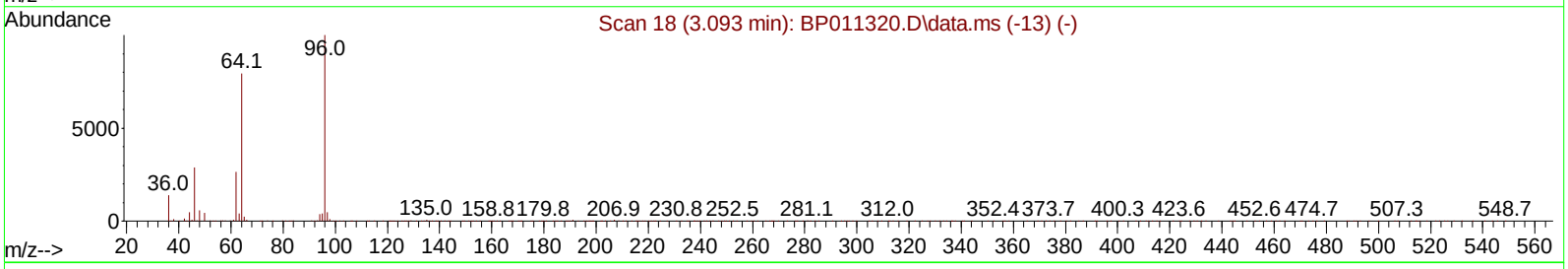

TIC: BP011335.D\data.ms

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

3.093min (+ 0.006) 3.55 ng/uL m

response 13651

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	81.81#
34.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

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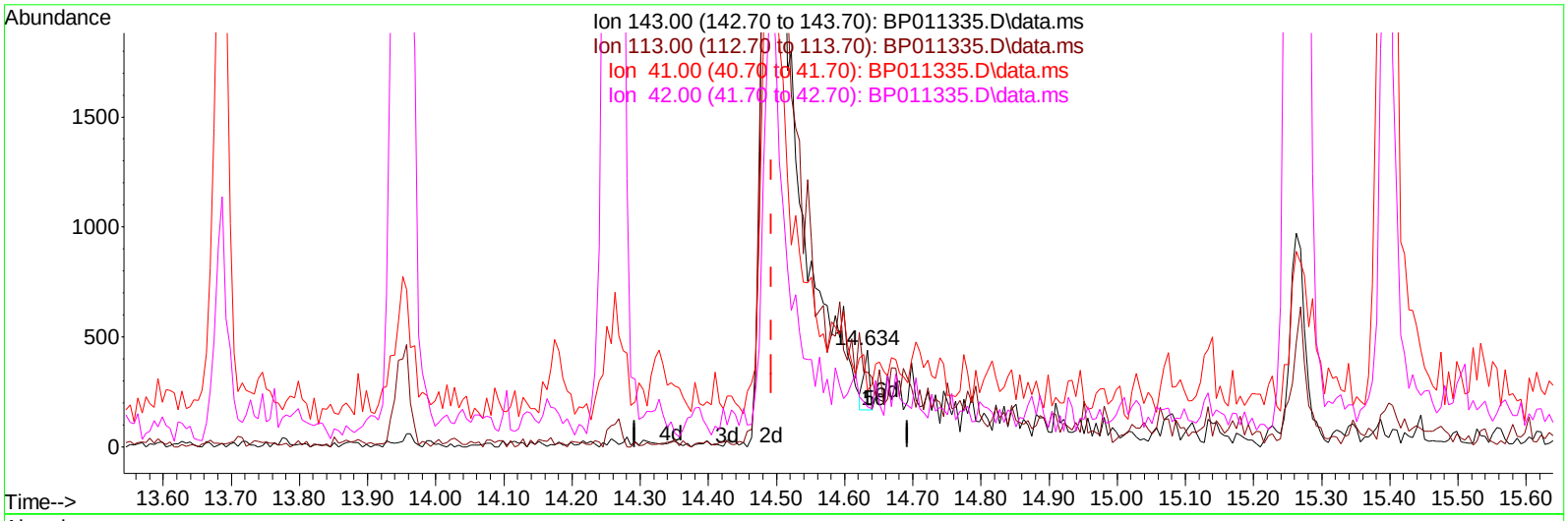

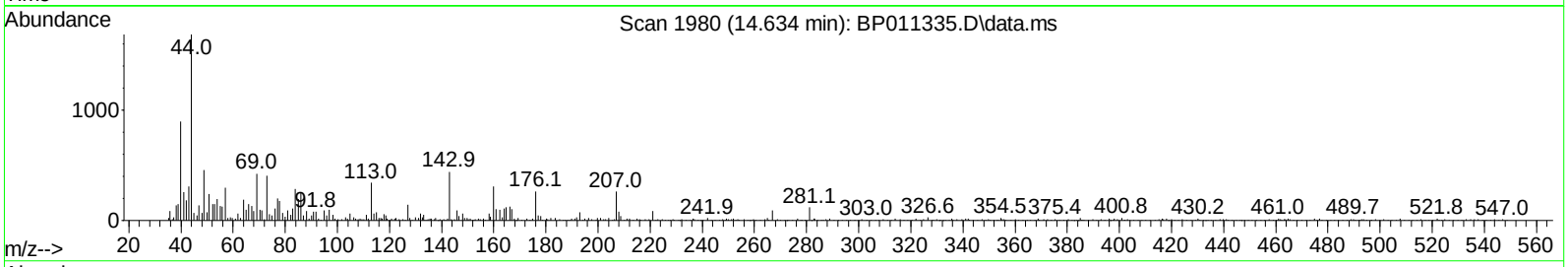

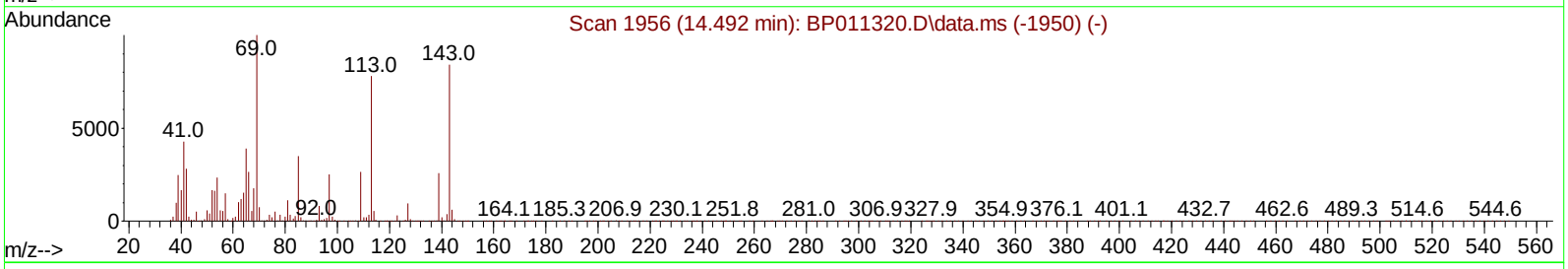

TIC: BP011335.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.634min (+ 0.141) 0.04 ng/ul

response 152

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	77.27
41.00	54.40	58.18
42.00	35.50	41.59

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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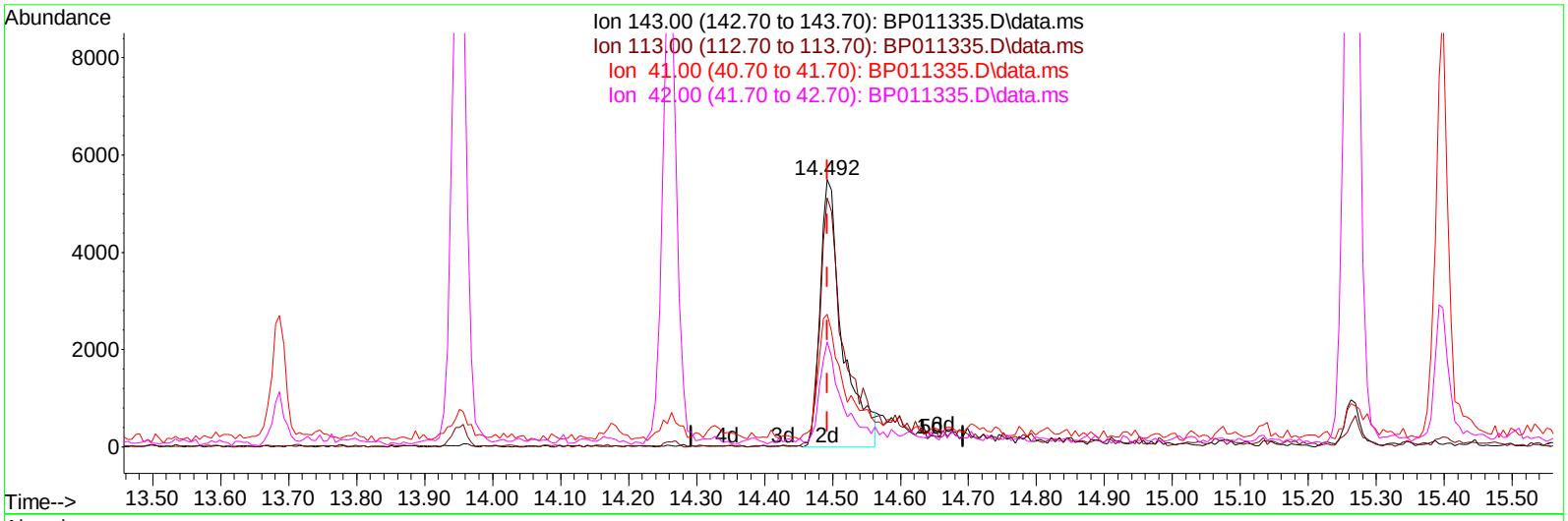

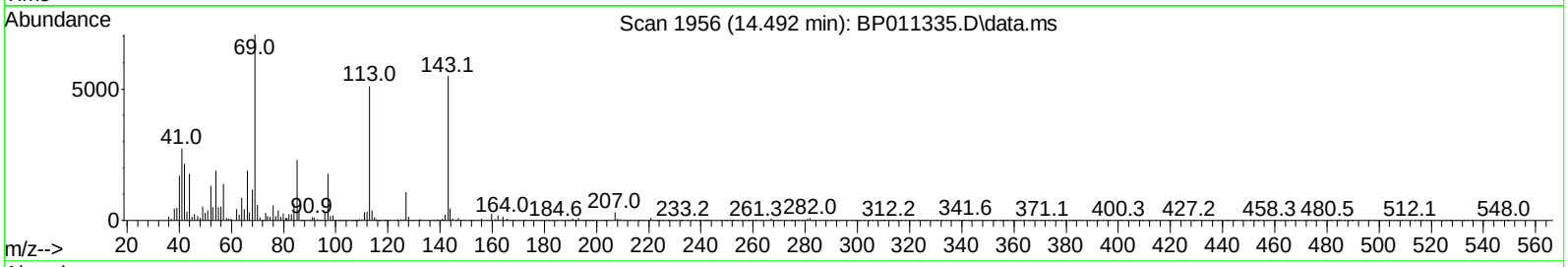

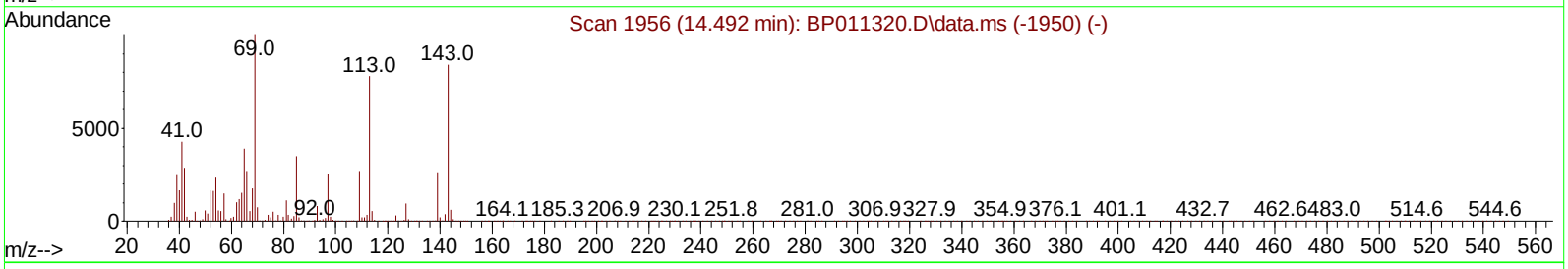

TIC: BP011335.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.492min (-0.000) 3.52 ng/ul m

response 12692

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	93.24#
41.00	54.40	49.65
42.00	35.50	39.31

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	117966	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	499156	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	284663	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	560569	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	519392	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	528419	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	13651m	3.547	ng/uL	0.00
4) Pyridine-d5	3.511	84	76661	7.898	ng/ul	0.01
7) Phenol-d5	6.775	99	64334	6.162	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	226987	30.665	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	196714	24.015	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	125115	15.263	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	123216	36.271	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	112270	36.097	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	191333	30.462	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	282125	25.693	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	763986	38.680	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	861218	36.223	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	12692m	3.518	ng/ul	0.00
60) Fluorene-d10	15.263	176	621824	36.813	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	86410	33.186	ng/ul	0.00
73) Anthracene-d10	17.133	188	1033520	41.019	ng/ul	0.00
81) Pyrene-d10	19.392	212	1169340	40.864	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1101010	41.316	ng/ul	0.00
Target Compounds						Qvalue

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

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