Quantitation Report (Qedit)

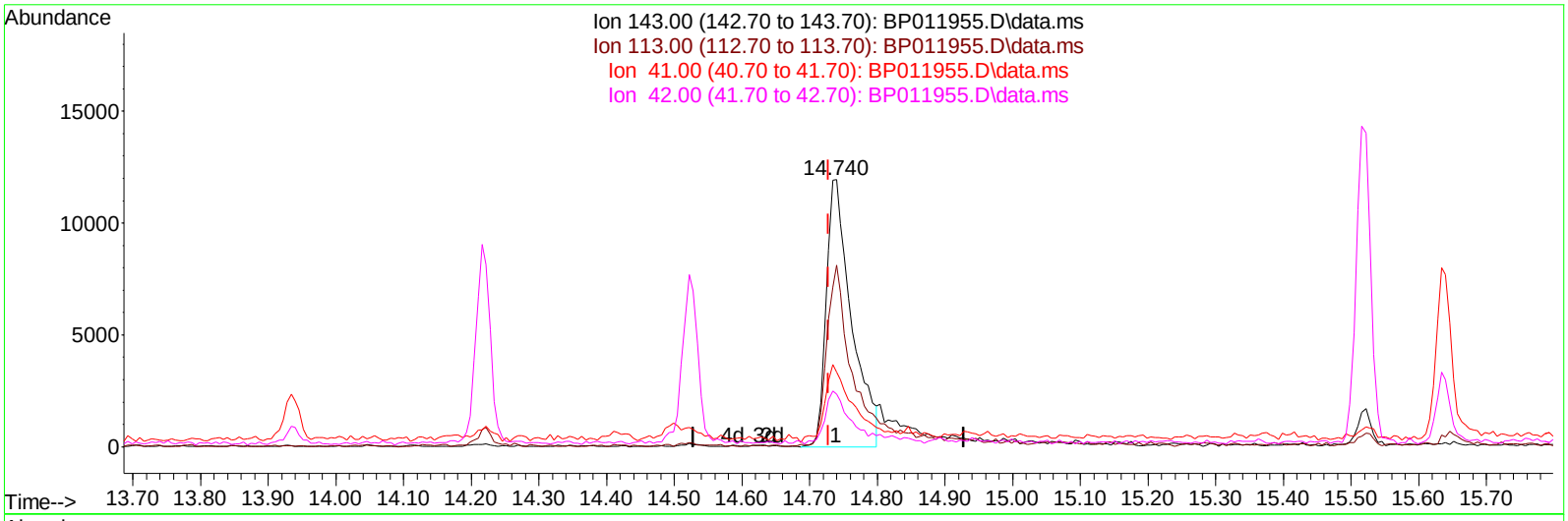
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP011955.D
Acq On : 06 Oct 2022 05:56
Operator : CG/JU
Sample : N4956-03
Misc :
ALS Vial : 35 Sample Multiplier: 1

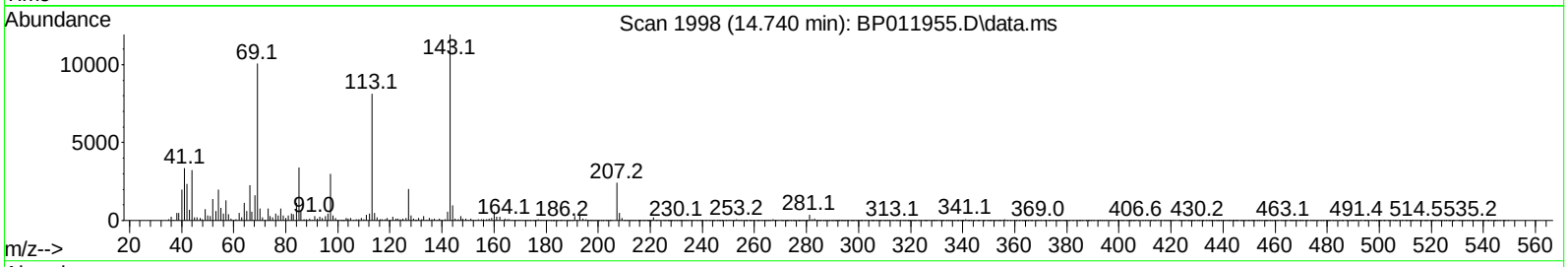
Quant Time: Oct 06 06:36:42 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 05 23:18:55 2022
Response via : Initial Calibration

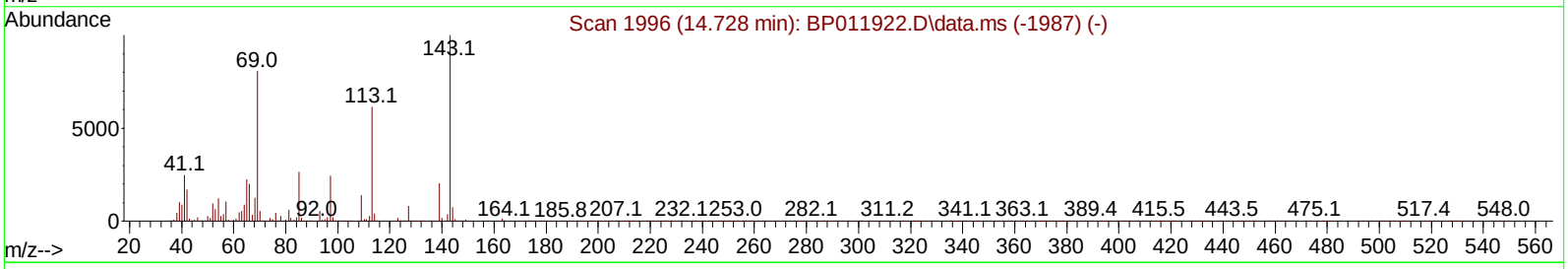
Instrument :
BNA_P
ClientSampleId :
CB8R8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/06/2022
Supervised By :mohammad ahmed 10/10/2022







TIC: BP011955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.64 ng/ul

response 31069

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	68.14
41.00	24.20	28.01
42.00	16.40	19.75#

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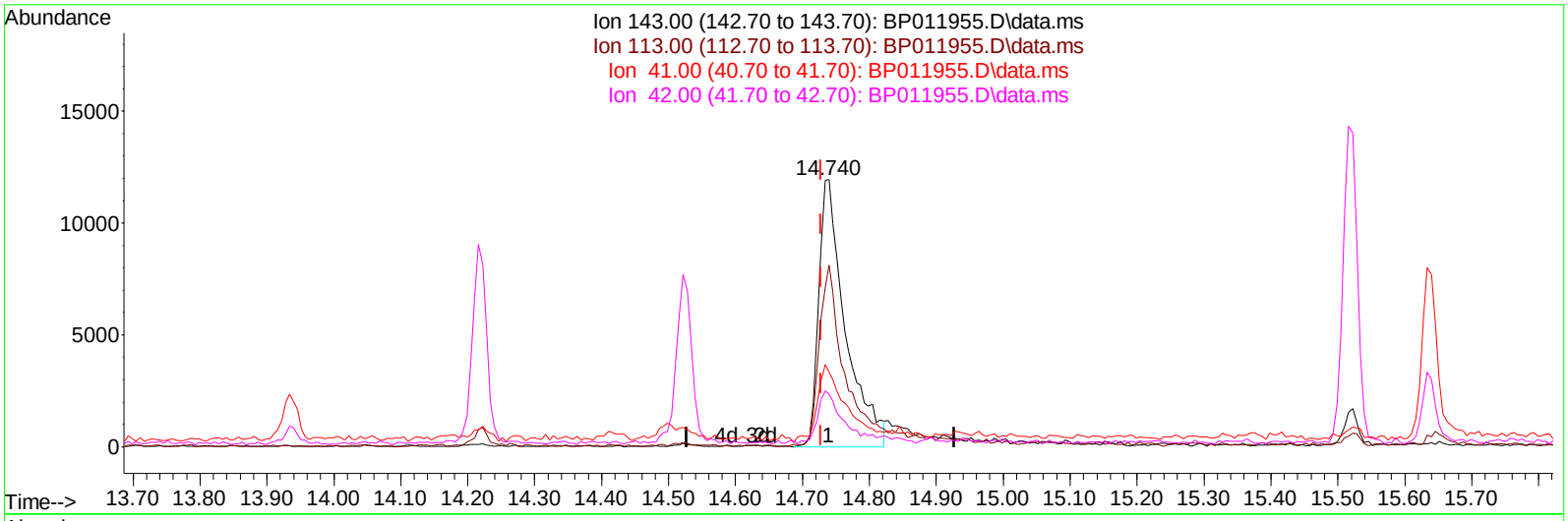

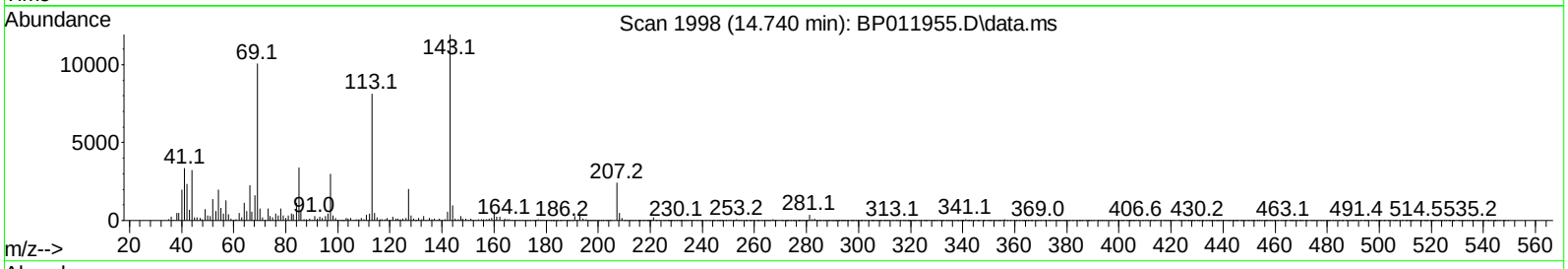

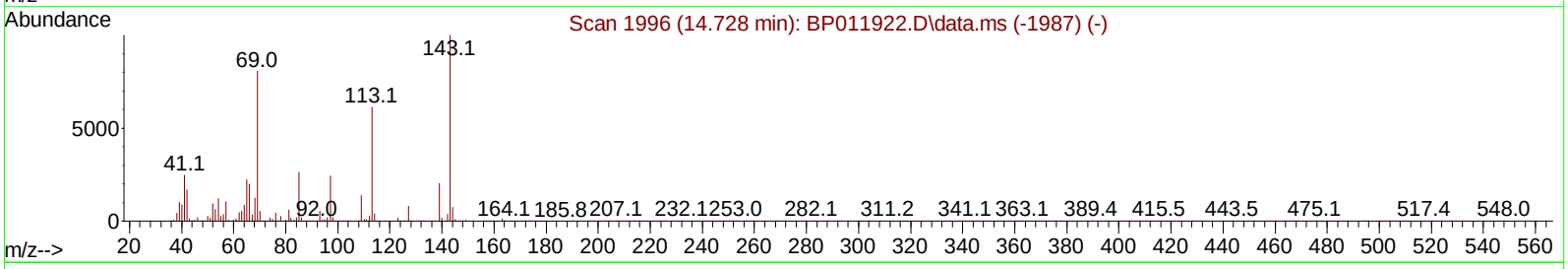

TIC: BP011955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.012) 3.87 ng/ul m

response 32966

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	68.14
41.00	24.20	28.01
42.00	16.40	19.75#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	215278	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	931676	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	575963	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1223562	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1160020	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1135764	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	25980	5.322	ng/uL	0.00
4) Pyridine-d5	3.728	84	111292	7.784	ng/ul	0.00
7) Phenol-d5	7.040	99	118531	6.378	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	363329	35.269	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	384165	26.185	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	222820	14.503	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	250404	35.286	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	241297	31.781	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	426356	30.360	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	438615	20.487	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1480710	35.951	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1691336	36.021	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	32966m	3.865	ng/ul	0.01
60) Fluorene-d10	15.522	176	1352187	37.803	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	184911	24.968	ng/ul	0.00
73) Anthracene-d10	17.381	188	2138036	38.360	ng/ul	0.00
81) Pyrene-d10	19.616	212	2494586	42.462	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2337955	41.121	ng/ul	0.00

Target Compounds Qvalue

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

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