Quantitation Report (Qedit)

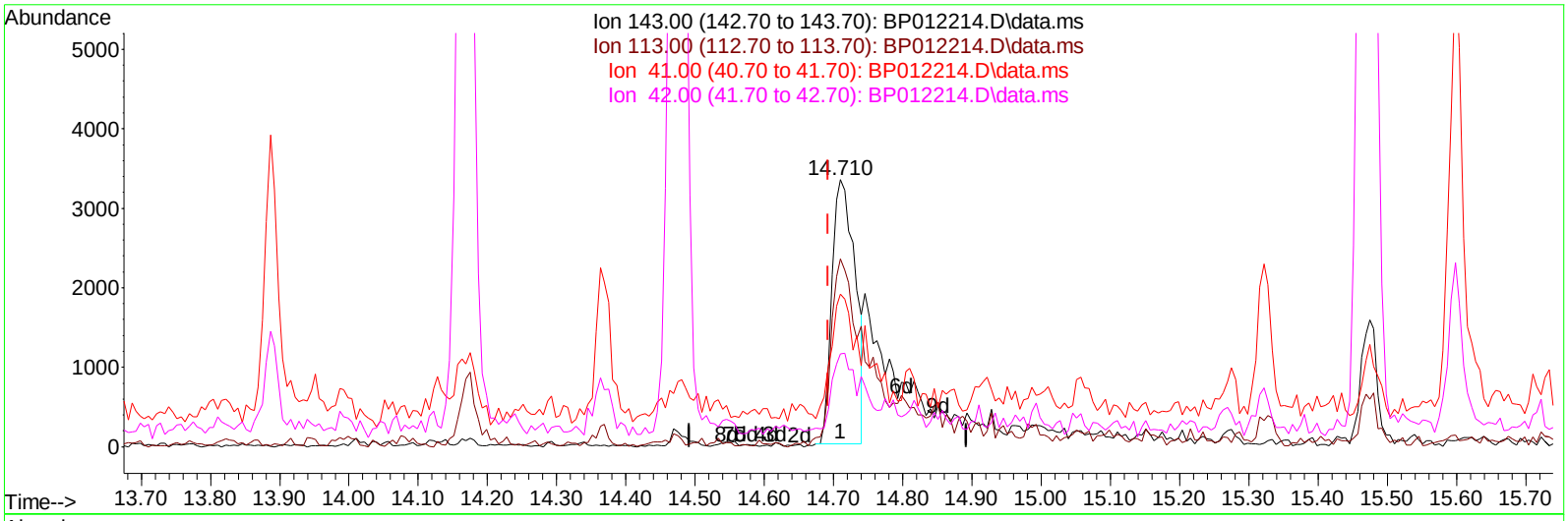
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
Data File : BP012214.D
Acq On : 20 Oct 2022 15:23
Operator : CG/JU
Sample : N5103-10
Misc :
ALS Vial : 9 Sample Multiplier: 1

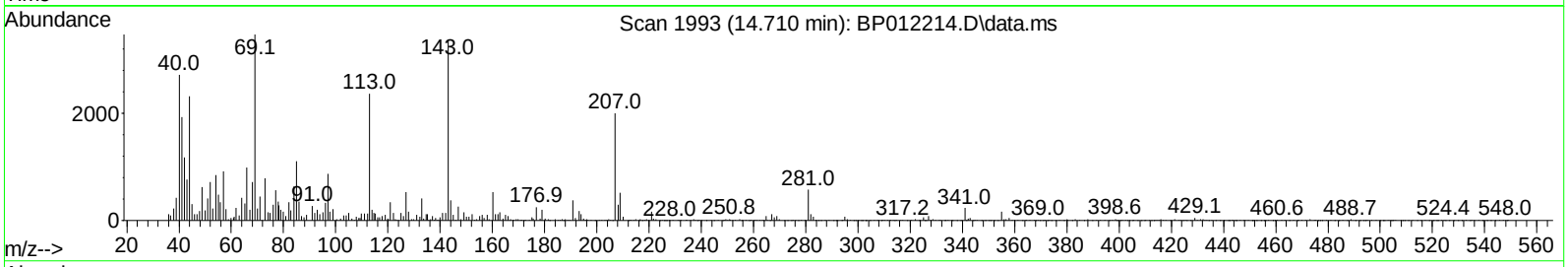
Quant Time: Oct 20 23:23:20 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 20 00:51:41 2022
Response via : Initial Calibration

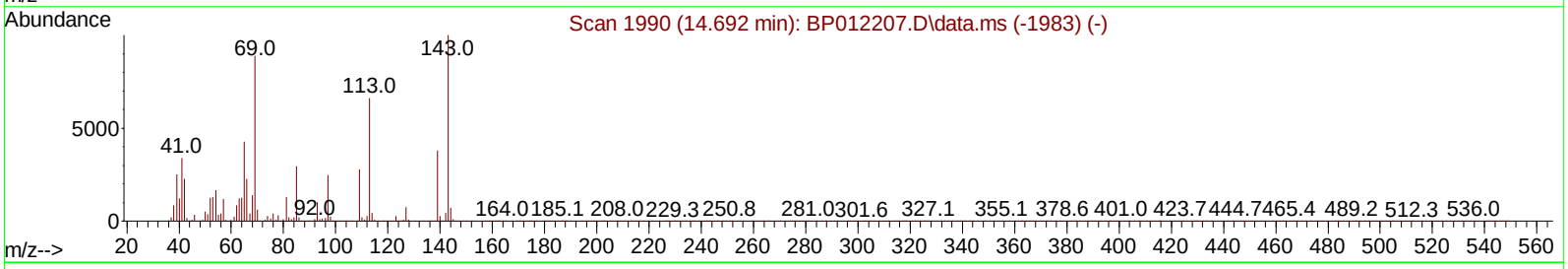
Instrument :
BNA_P
ClientSampleId :
C0BB3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 10/21/2022
Supervised By :mohammad ahmed 10/24/2022







TIC: BP012214.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.018) 0.70 ng/ul

response 7652

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.20	70.31
41.00	35.10	57.24#
42.00	23.90	34.90#

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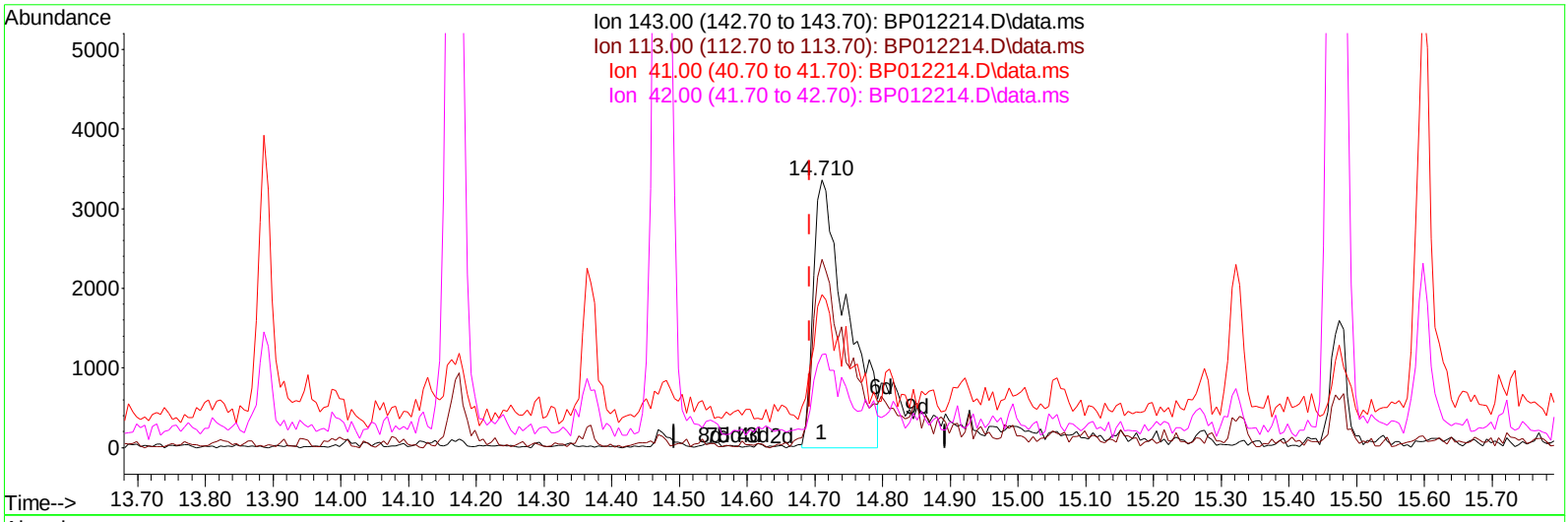

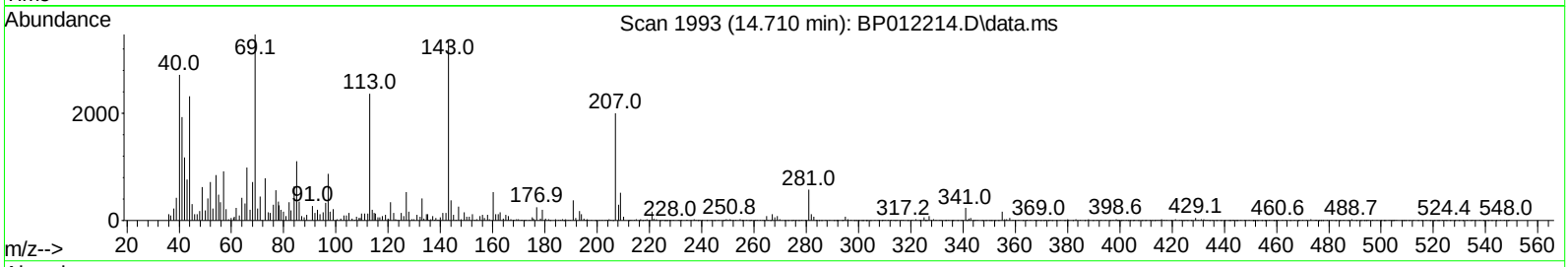

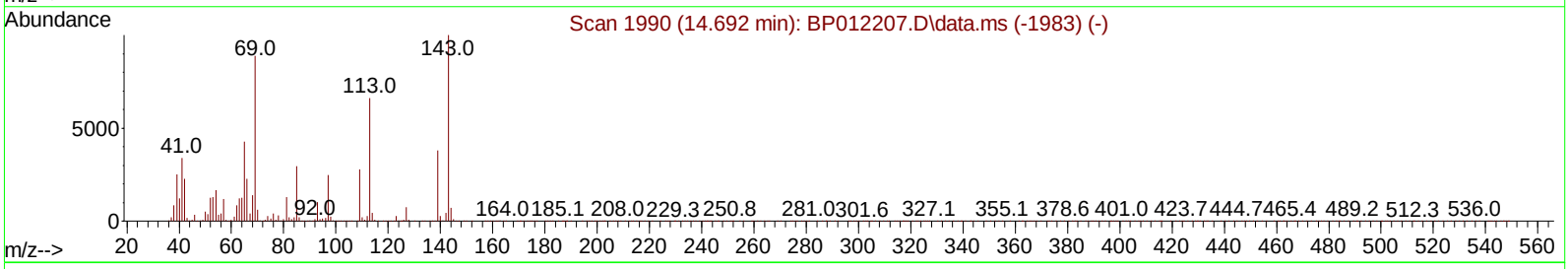

TIC: BP012214.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.018) 1.07 ng/ul m

response 11637

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.20	70.31
41.00	35.10	57.24#
42.00	23.90	34.90#

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	363593	20.000	ng/ul	0.00
20) Naphthalene-d8	10.628	136	1456165	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	838178	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.234	188	1640820	20.000	ng/ul	0.00
79) Chrysene-d12	21.327	240	1356628	20.000	ng/ul	0.00
88) Perylene-d12	23.739	264	1366248	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	36837	4.100	ng/uL	0.00
4) Pyridine-d5	3.687	84	197620	7.634	ng/ul	0.00
7) Phenol-d5	6.999	99	88084	2.806	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	505566	26.558	ng/ul	0.00
11) 2-Chlorophenol-d4	7.358	132	283497	11.865	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	187971	7.705	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	300495	27.879	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	169214	14.551	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	302492	13.867	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	422231	13.624	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	1680228	28.818	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	1895150	28.476	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	11637m	1.065	ng/ul	0.02
60) Fluorene-d10	15.475	176	1472581	29.358	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	91802	9.835	ng/ul	0.00
73) Anthracene-d10	17.334	188	2245201	31.217	ng/ul	0.00
81) Pyrene-d10	19.575	212	2393107	30.983	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264	2147542	31.344	ng/ul	0.00

Target Compounds Qvalue

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

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