Quantitation Report (Qedit)

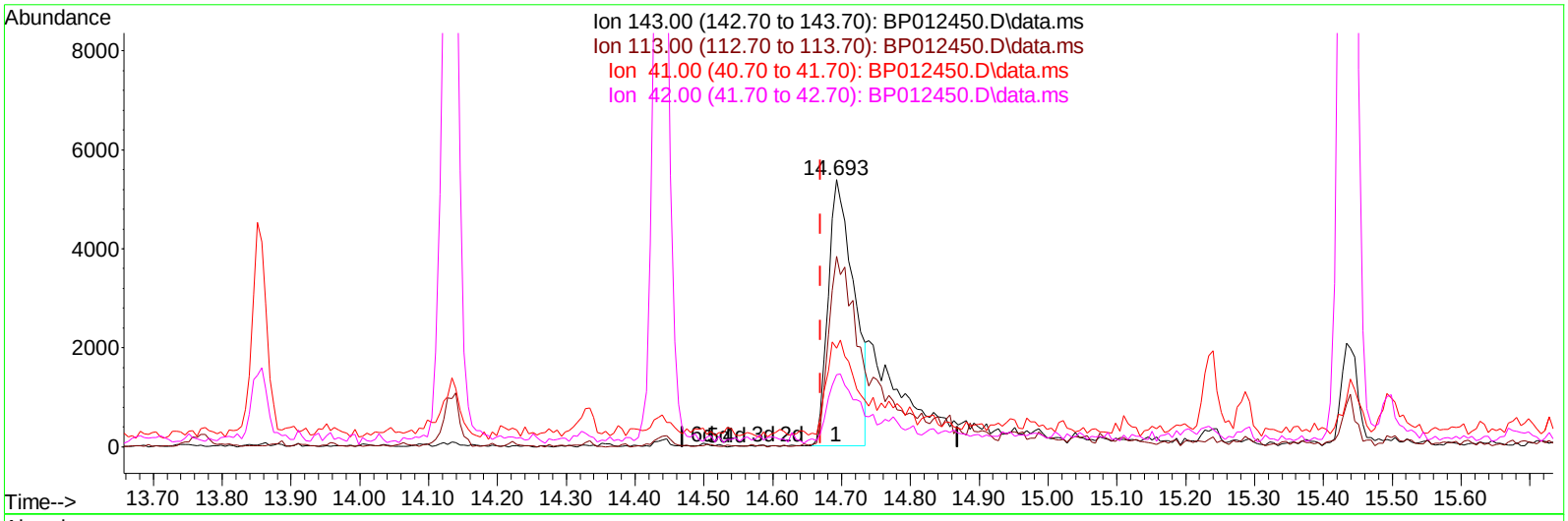
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012450.D
Acq On : 02 Nov 2022 13:01
Operator : CG/JU
Sample : N5180-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

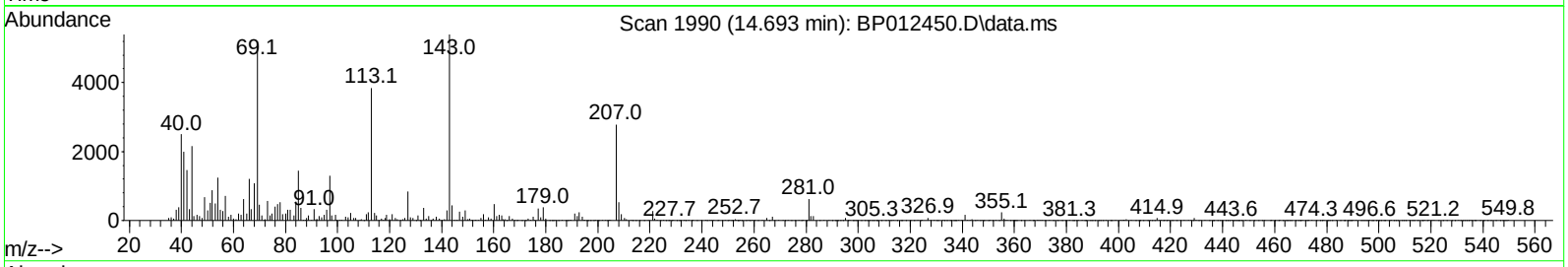
Quant Time: Nov 02 23:13:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 01 01:03:58 2022
Response via : Initial Calibration

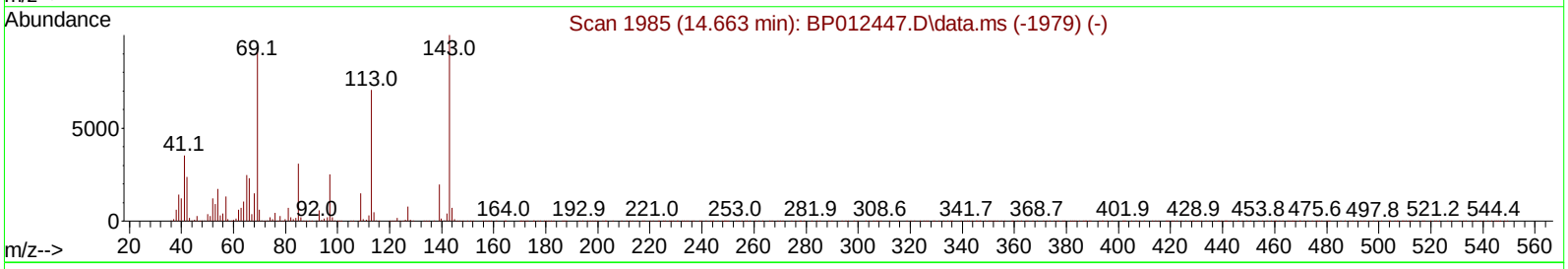
Instrument :
BNA_P
ClientSampleId :
C0KX2

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2022
Supervised By :mohammad ahmed 11/03/2022







TIC: BP012450.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.024) 1.23 ng/ul

response 13889

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	71.23
41.00	34.50	36.79
42.00	23.60	27.08

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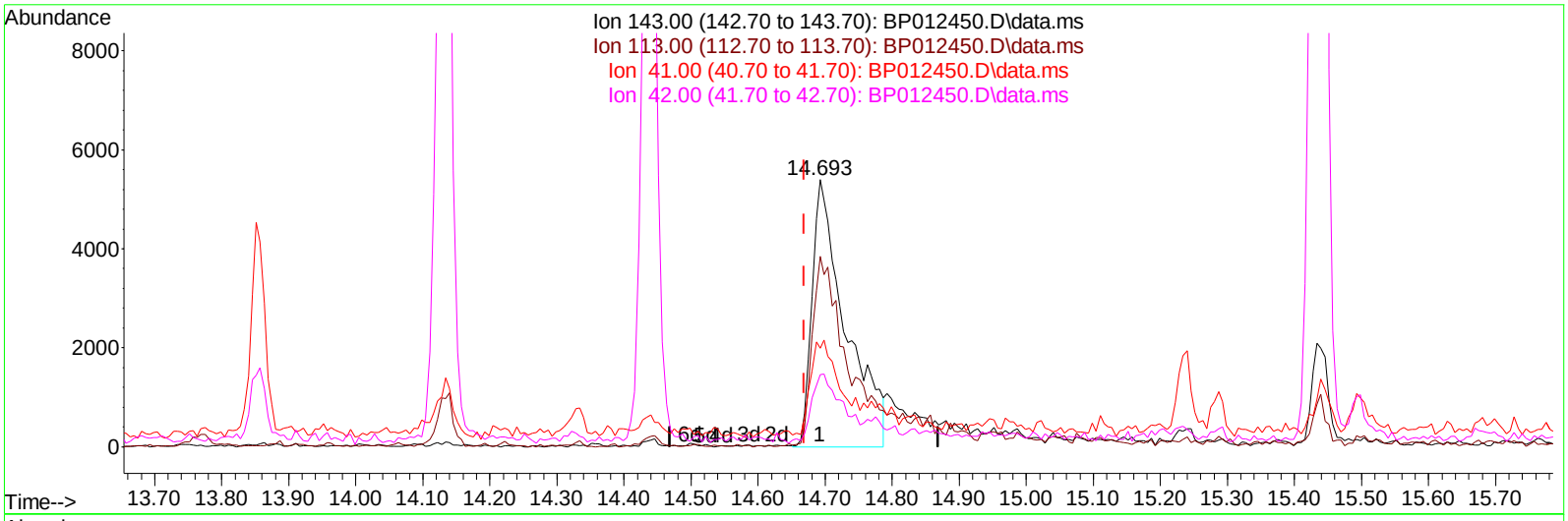

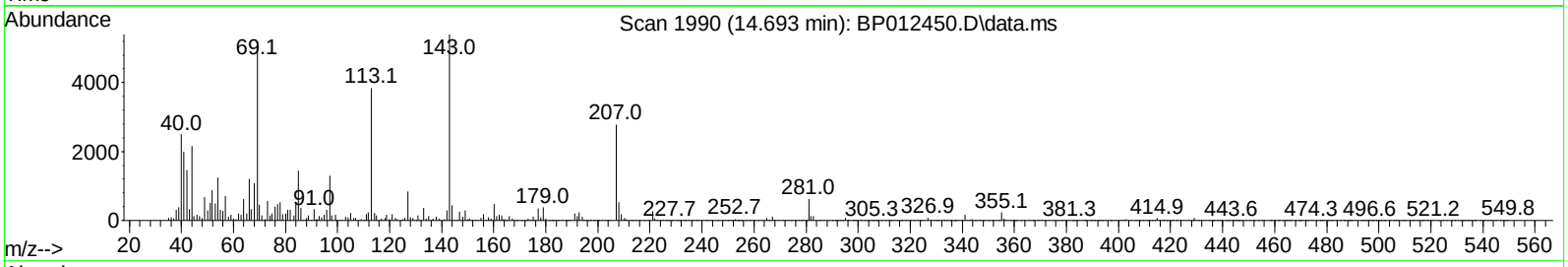

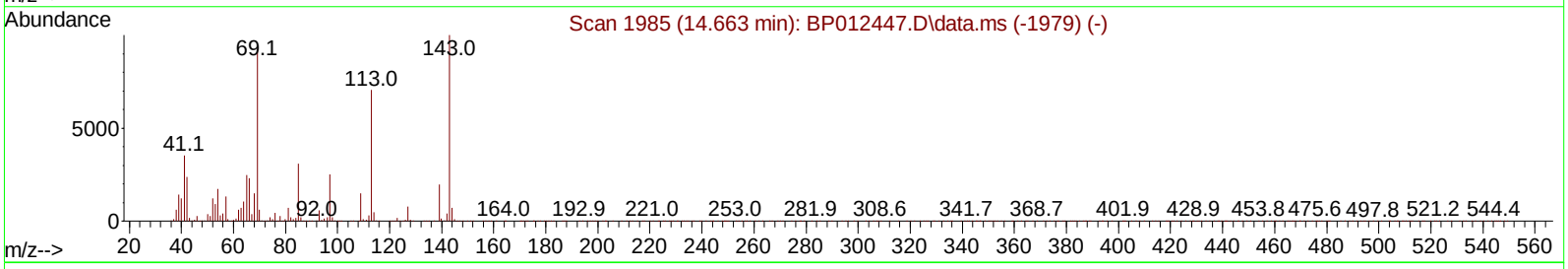

TIC: BP012450.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.024) 1.66 ng/ul m

response 18765

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	71.23
41.00	34.50	36.79
42.00	23.60	27.08

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	281172	20.000	ng/ul	0.00
20) Naphthalene-d8	10.587	136	1271556	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.440	164	875049	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	1959738	20.000	ng/ul	-0.01
79) Chrysene-d12	21.298	240	1865278	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1703507	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23316	3.339	ng/uL	0.00
4) Pyridine-d5	3.664	84	134763	6.718	ng/ul	0.00
7) Phenol-d5	6.964	99	162388	6.439	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	458347	30.445	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.322	132	186233	10.112	ng/ul	0.00
15) 4-Methylphenol-d8	8.511	113	208077	10.432	ng/ul	0.00
21) Nitrobenzene-d5	8.958	128	314286	38.115	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	70412	8.227	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	211739	11.416	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	638849	23.132	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2274354	37.968	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2420223	35.355	ng/ul	0.00
54) 4-Nitrophenol-d4	14.693	143	18765m	1.655	ng/ul	0.02
60) Fluorene-d10	15.440	176	1982120	38.648	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	2798	0.274	ng/ul	0.01
73) Anthracene-d10	17.304	188	3444808	40.796	ng/ul	0.00
81) Pyrene-d10	19.545	212	3922675	41.887	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	3399570	40.843	ng/ul	-0.01
Target Compounds						Qvalue

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

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