Quantitation Report (Qedit)

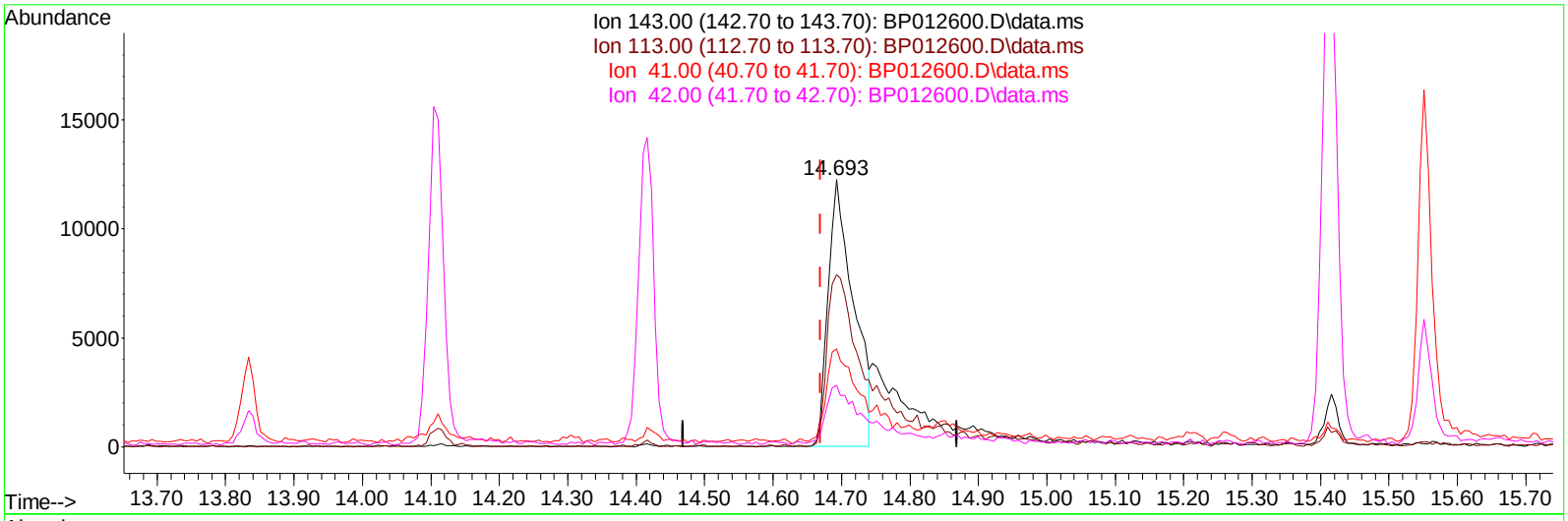
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
Data File : BP012600.D
Acq On : 10 Nov 2022 15:32
Operator : CG/JU
Sample : N5493-03
Misc :
ALS Vial : 8 Sample Multiplier: 1

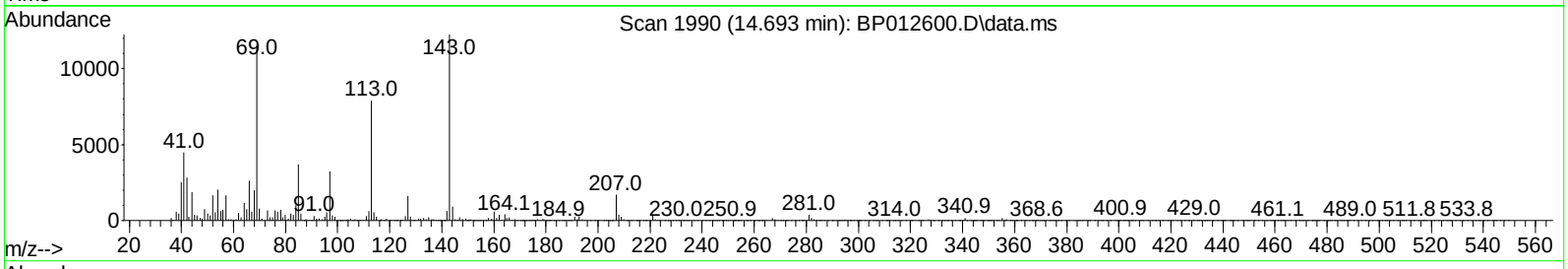
Quant Time: Nov 10 21:35:26 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 09 23:28:28 2022
Response via : Initial Calibration

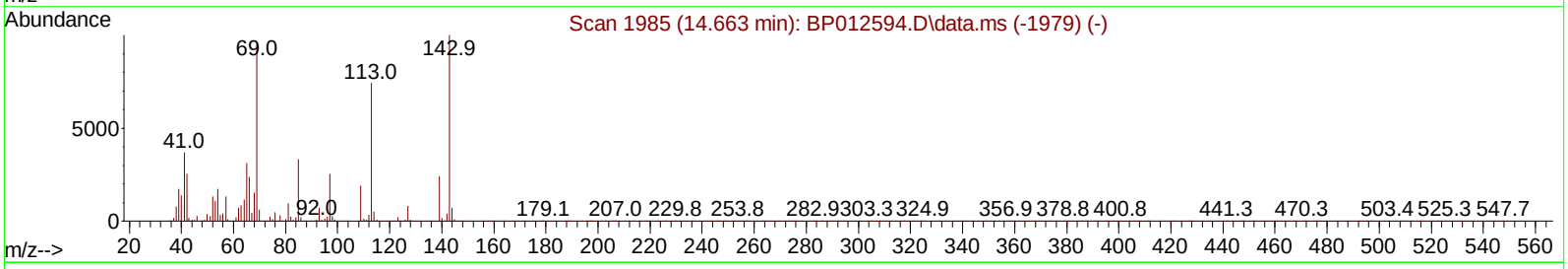
Instrument :
BNA_P
ClientSampleId :
H0AA4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/11/2022
Supervised By :mohammad ahmed 11/14/2022







TIC: BP012600.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.024) 3.42 ng/ul

response 31546

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	64.46
41.00	35.20	36.62
42.00	24.00	23.02

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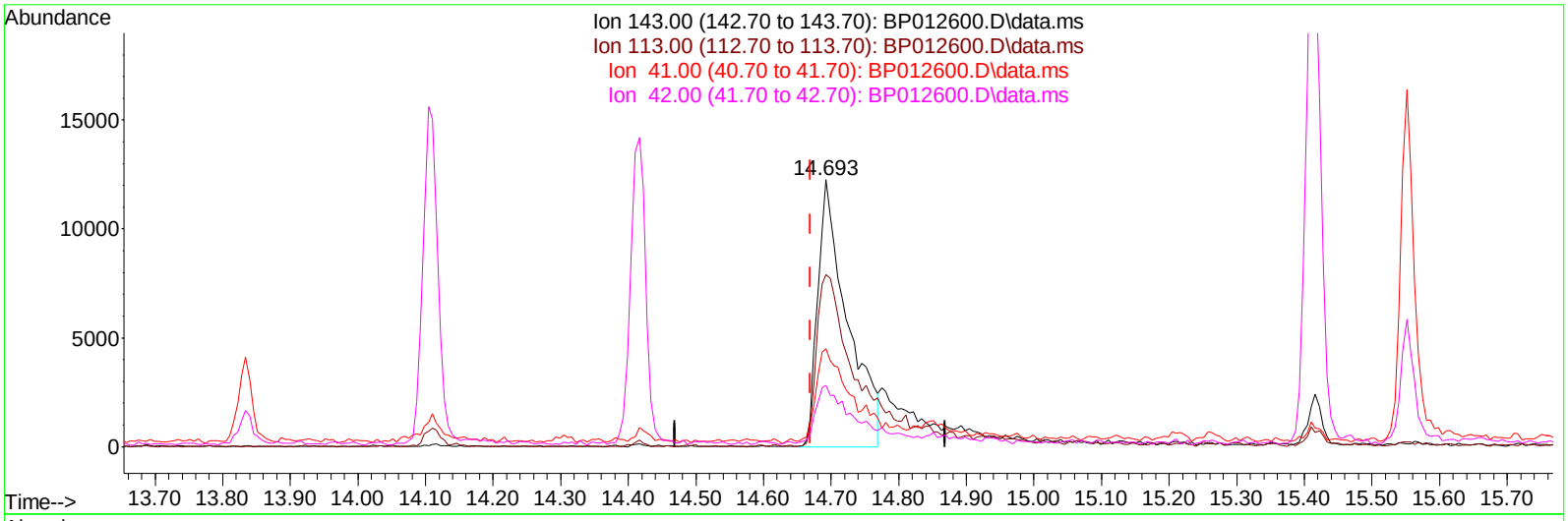

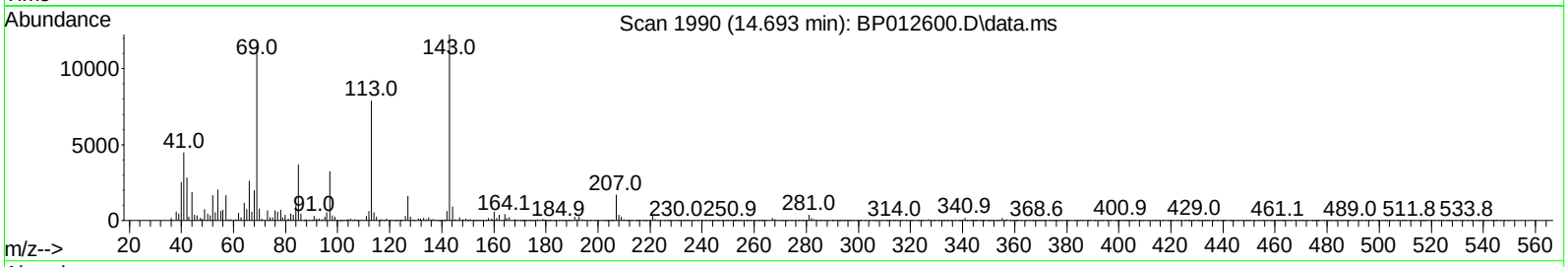

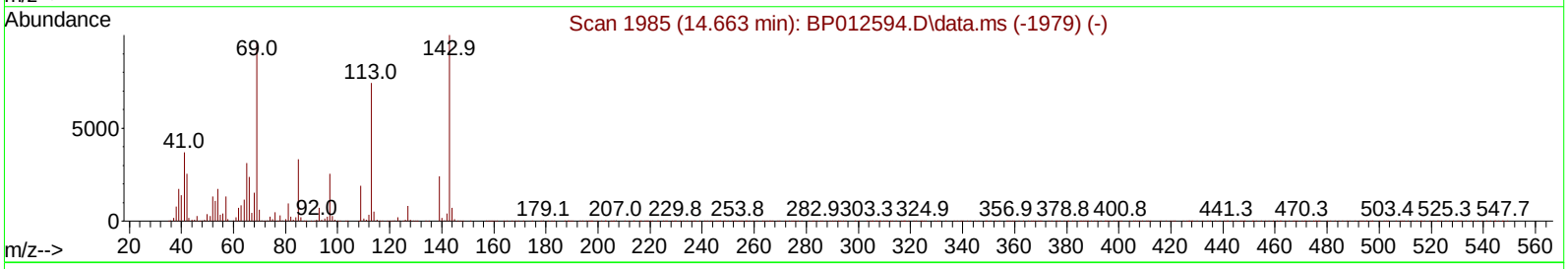

TIC: BP012600.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.693min (+ 0.024) 4.05 ng/ul m

response 37280

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	64.46
41.00	35.20	36.62
42.00	24.00	23.02

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.764	152	274445	20.000	ng/ul	0.00
20) Naphthalene-d8	10.558	136	1271161	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	849089	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1873982	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1936950	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	1869768	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	29238	4.342	ng/uL	0.00
4) Pyridine-d5	3.652	84	154605	8.125	ng/ul	0.01
7) Phenol-d5	6.952	99	143912	6.220	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	483995	34.964	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	448686	25.202	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	283130	15.471	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	315789	33.575	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	325460	31.505	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	512189	27.521	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	736668	28.101	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1998898	34.154	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	2253928	34.043	ng/ul	0.00
54) 4-Nitrophenol-d4	14.693	143	37280m	4.046	ng/ul	0.02
60) Fluorene-d10	15.416	176	1787668	36.024	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	306561	29.577	ng/ul	0.00
73) Anthracene-d10	17.281	188	2978281	36.597	ng/ul	0.00
81) Pyrene-d10	19.522	212	3466018	31.932	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.498	264	3365355	36.511	ng/ul	0.00
Target Compounds						Qvalue

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

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