Quantitation Report (Qedit)

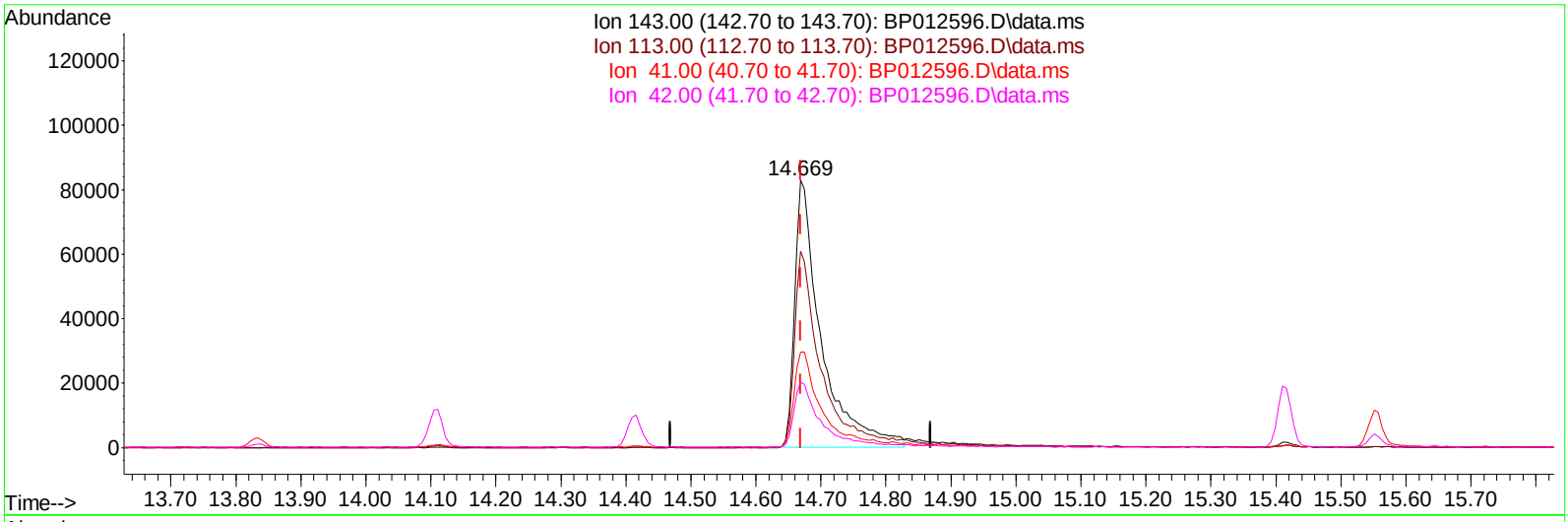
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
Data File : BP012596.D
Acq On : 10 Nov 2022 13:14
Operator : CG/JU
Sample : PB148942BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

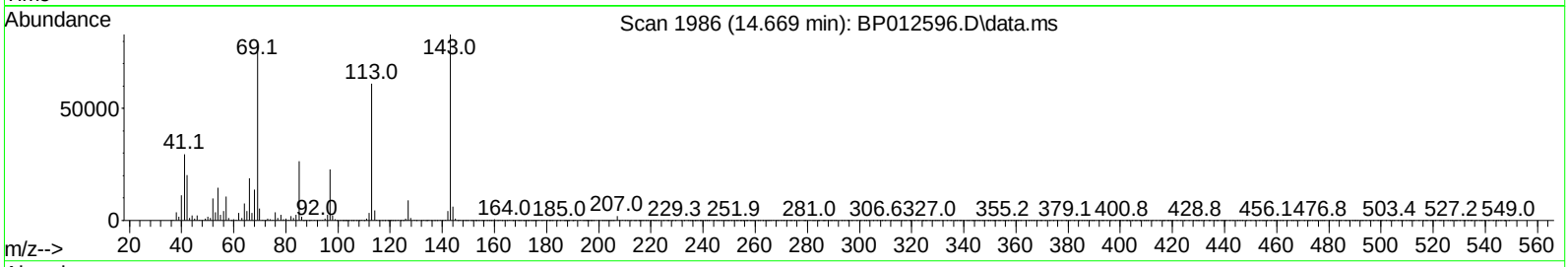
Quant Time: Nov 10 21:34:44 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 09 23:28:28 2022
Response via : Initial Calibration

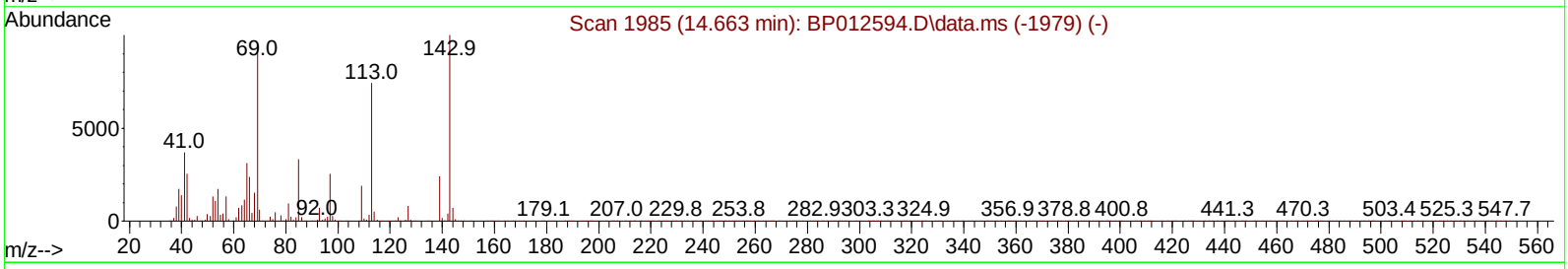
Instrument :
BNA_P
ClientSampleId :
SBLK942

Manual IntegrationsAPPROVED

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







TIC: BP012596.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 36.13 ng/ul

response 234978

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.51
41.00	35.20	35.69
42.00	24.00	24.44

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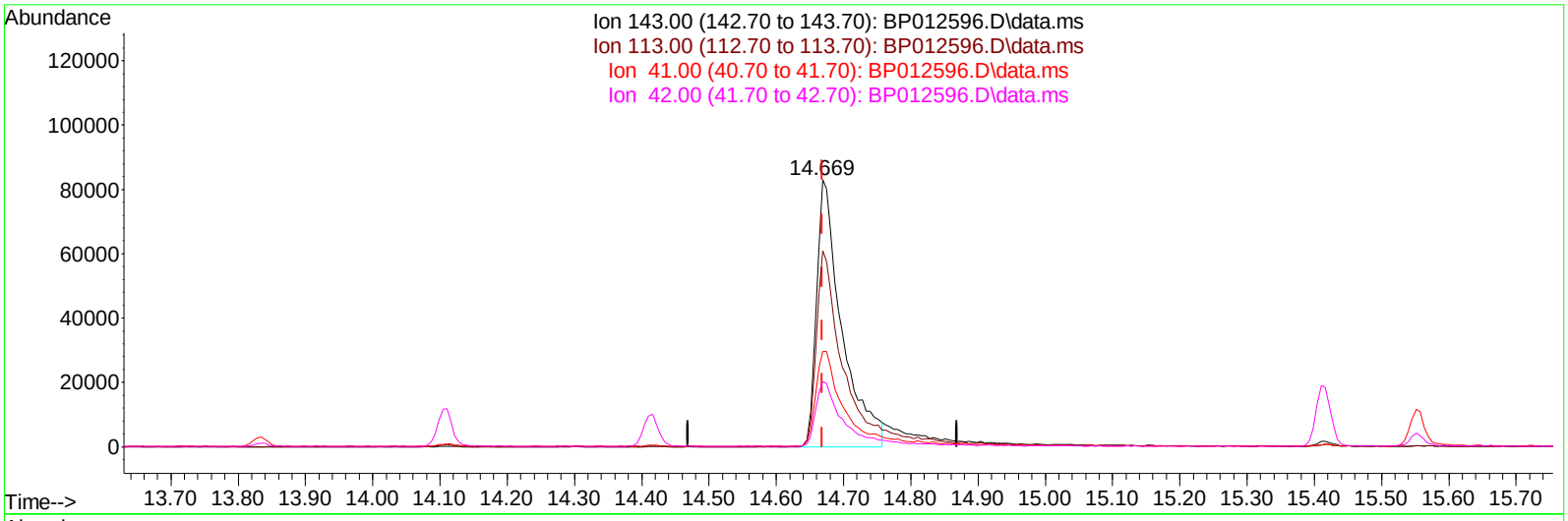

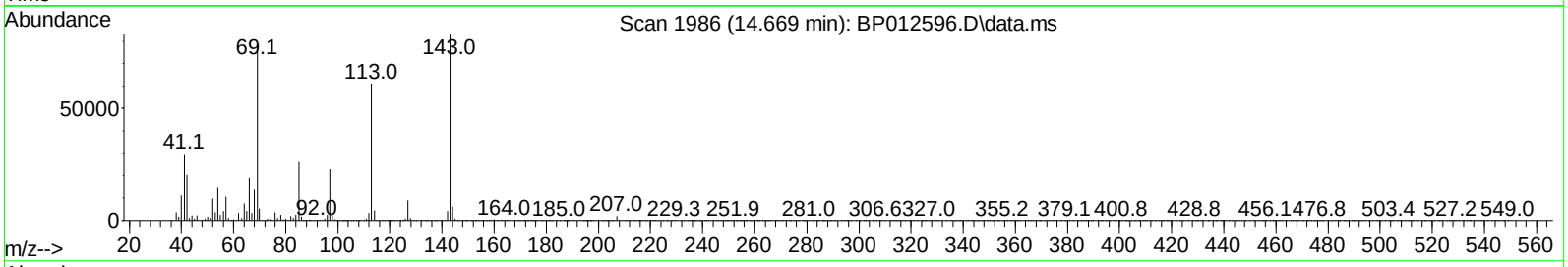

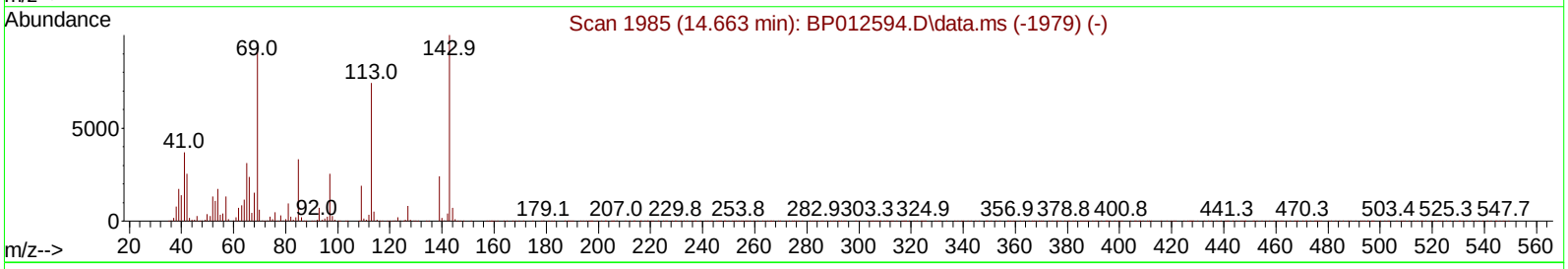

TIC: BP012596.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 33.30 ng/ul m

response 216559

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.51
41.00	35.20	35.69
42.00	24.00	24.44

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	199348	20.000	ng/ul	0.00
20) Naphthalene-d8	10.557	136	902728	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	599265	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1326840	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1386689	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	1321620	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	35470	7.252	ng/uL	0.00
4) Pyridine-d5	3.646	84	190942	13.815	ng/ul	0.00
7) Phenol-d5	6.952	99	595960	35.463	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	379280	37.721	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	477653	36.935	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	477915	35.951	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	247268	37.019	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	267125	36.412	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	452125	34.208	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	331786	17.822	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1549940	37.524	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1717412	36.753	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	216559m	33.298	ng/ul	0.00
60) Fluorene-d10	15.416	176	1352049	38.604	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	234787	31.993	ng/ul	0.00
73) Anthracene-d10	17.281	188	2248371	39.021	ng/ul	0.00
81) Pyrene-d10	19.522	212	2629492	33.838	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.498	264	2595327	39.835	ng/ul	0.00
Target Compounds						Qvalue

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

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