Quantitation Report (Qedit)

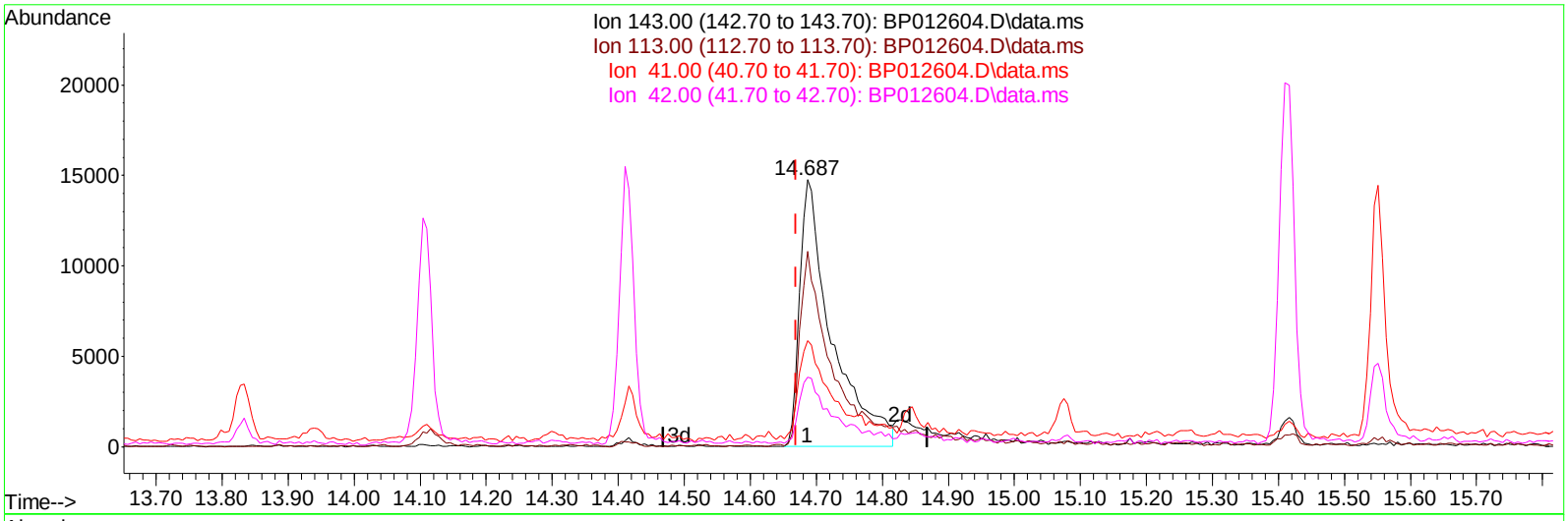
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
Data File : BP012604.D
Acq On : 10 Nov 2022 17:50
Operator : CG/JU
Sample : N5493-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

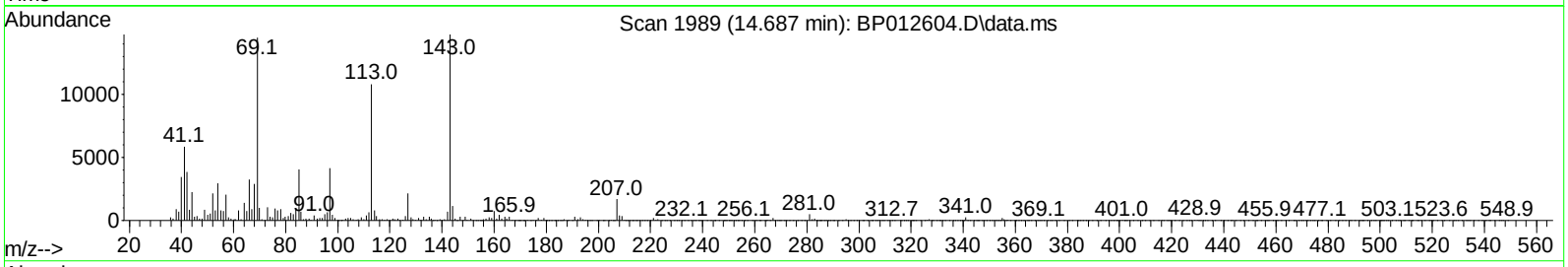
Quant Time: Nov 10 21:36:14 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 09 23:28:28 2022
Response via : Initial Calibration

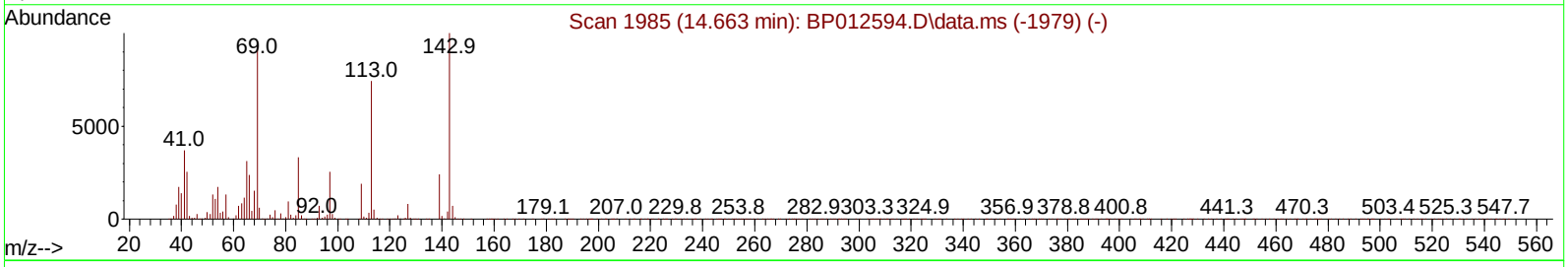
Instrument :
BNA_P
ClientSampleId :
H0AC5

Manual IntegrationsAPPROVED

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







TIC: BP012604.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.018) 5.08 ng/ul

response 49482

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.23
41.00	35.20	39.63
42.00	24.00	25.96

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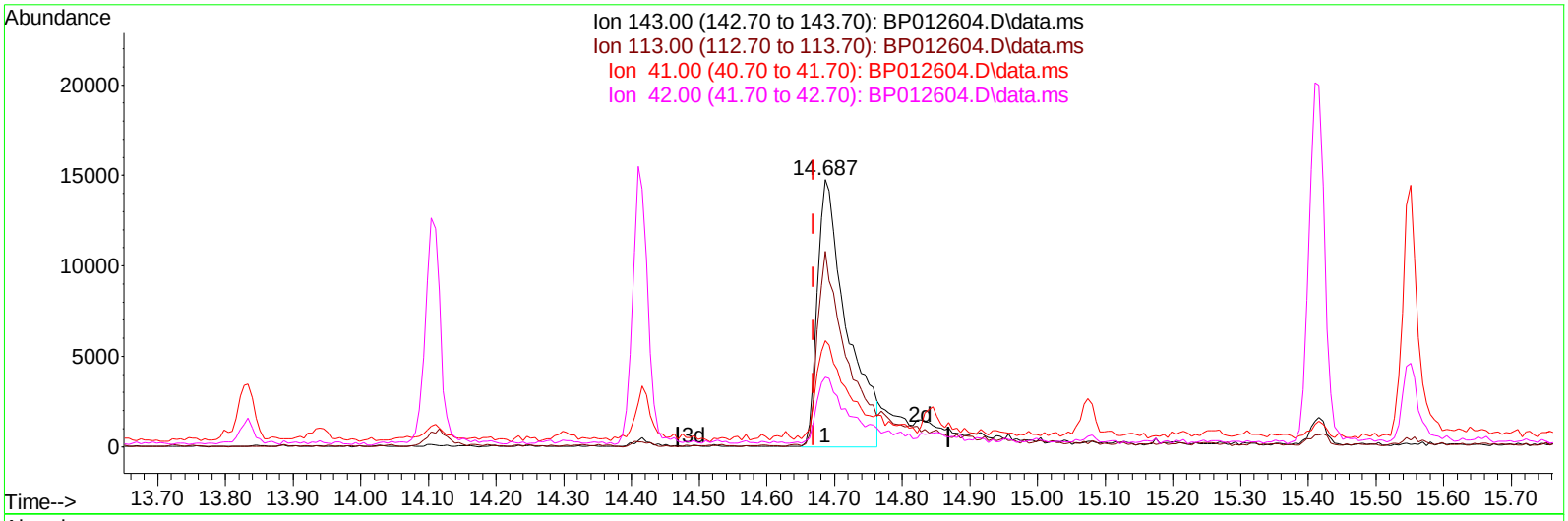

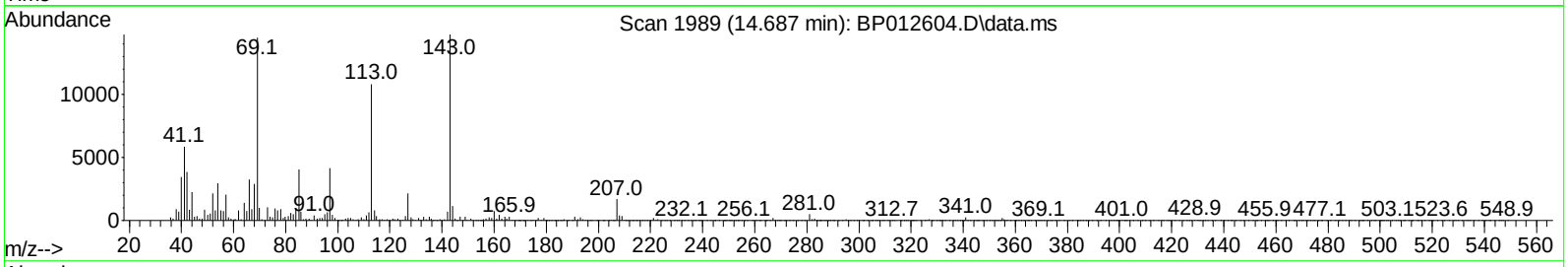

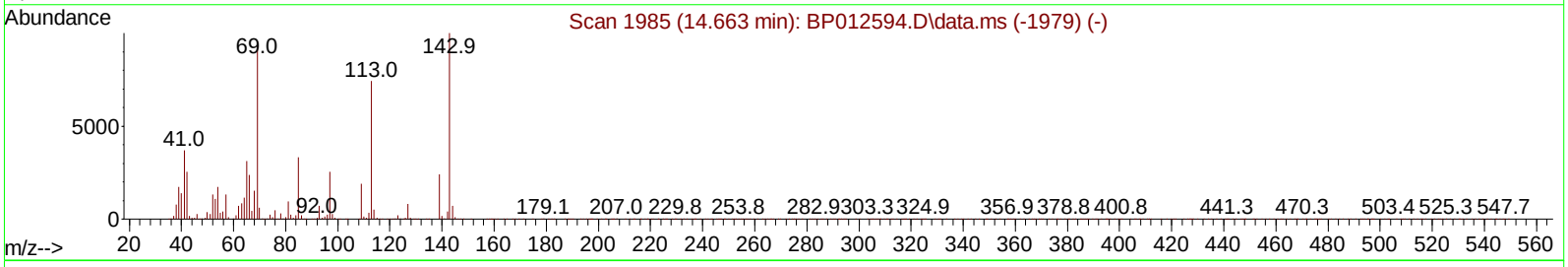

TIC: BP012604.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (+ 0.018) 4.55 ng/ul m

response 44340

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	71.10	73.23
41.00	35.20	39.63
42.00	24.00	25.96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	289035	20.000	ng/ul	0.00
20) Naphthalene-d8	10.557	136	1336762	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	897544	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.180	188	1966385	20.000	ng/ul	0.00
79) Chrysene-d12	21.274	240	2052037	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	2030105	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	27260	3.844	ng/uL	0.00
4) Pyridine-d5	3.652	84	110384	5.508	ng/ul	0.01
7) Phenol-d5	6.952	99	139170	5.712	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	404303	27.733	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	395355	21.085	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	269356	13.975	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	258152	26.100	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	273340	25.161	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	445600	22.768	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	500890	18.169	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1784810	28.850	ng/ul	0.00
49) Acenaphthylene-d8	14.104	160	1889479	26.998	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	44340m	4.552	ng/ul	0.02
60) Fluorene-d10	15.410	176	1548484	29.519	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	279784	25.725	ng/ul	0.00
73) Anthracene-d10	17.281	188	2736422	32.045	ng/ul	0.00
81) Pyrene-d10	19.522	212	3261411	28.362	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	3228798	32.263	ng/ul	0.00
Target Compounds						Qvalue

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

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