Quantitation Report (Qedit)

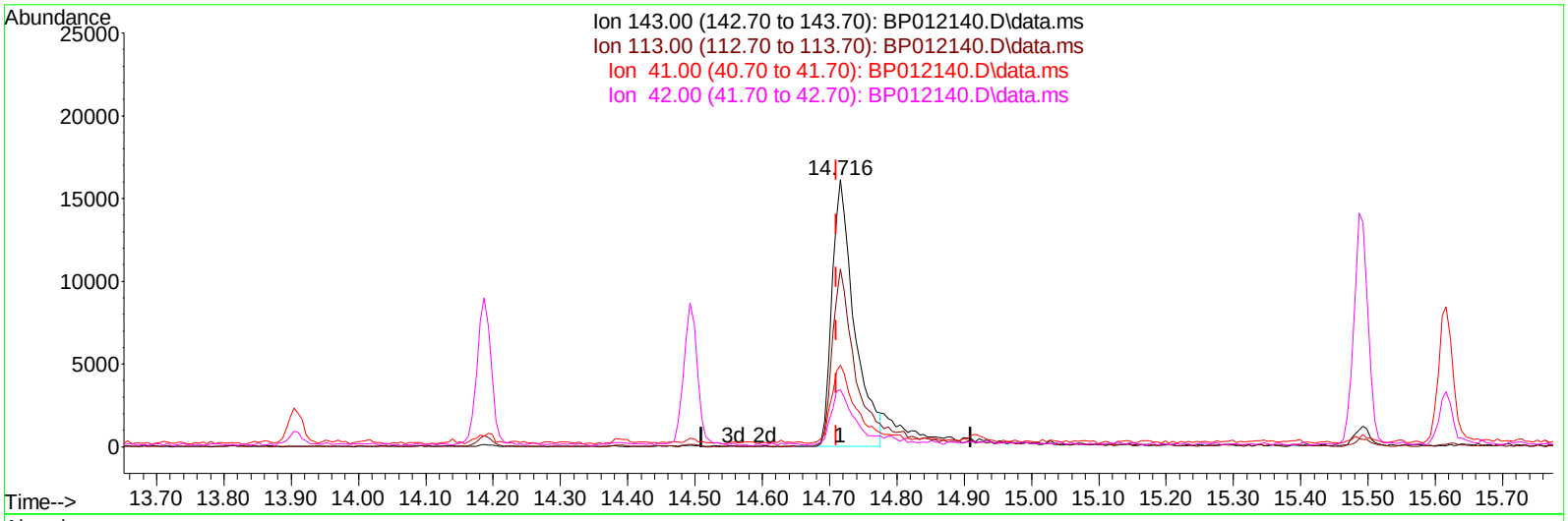
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
Data File : BP012140.D
Acq On : 14 Oct 2022 16:25
Operator : CG/JU
Sample : N4982-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

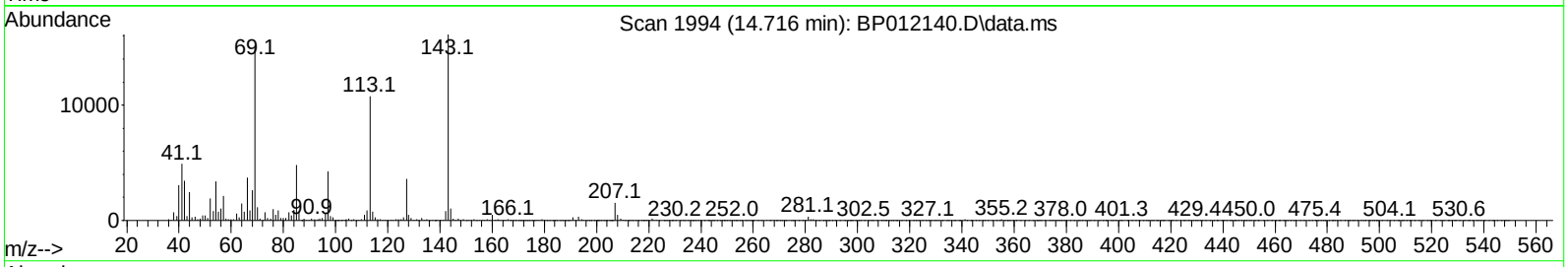
Quant Time: Oct 14 23:04:03 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 12 23:03:21 2022
Response via : Initial Calibration

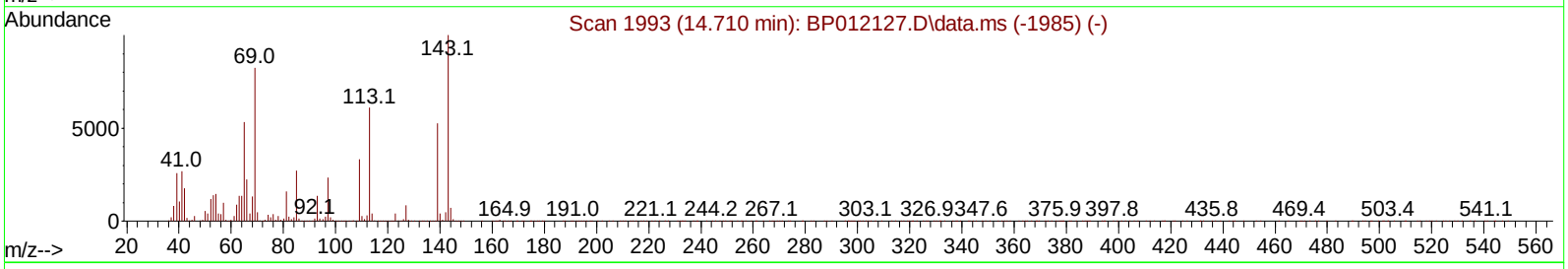
Instrument :
BNA_P
ClientSampleId :
C1BC0

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/15/2022
Supervised By :mohammad ahmed 10/16/2022







TIC: BP012140.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 5.16 ng/ul

response 37279

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.80	66.67
41.00	26.60	30.46
42.00	17.60	21.41#

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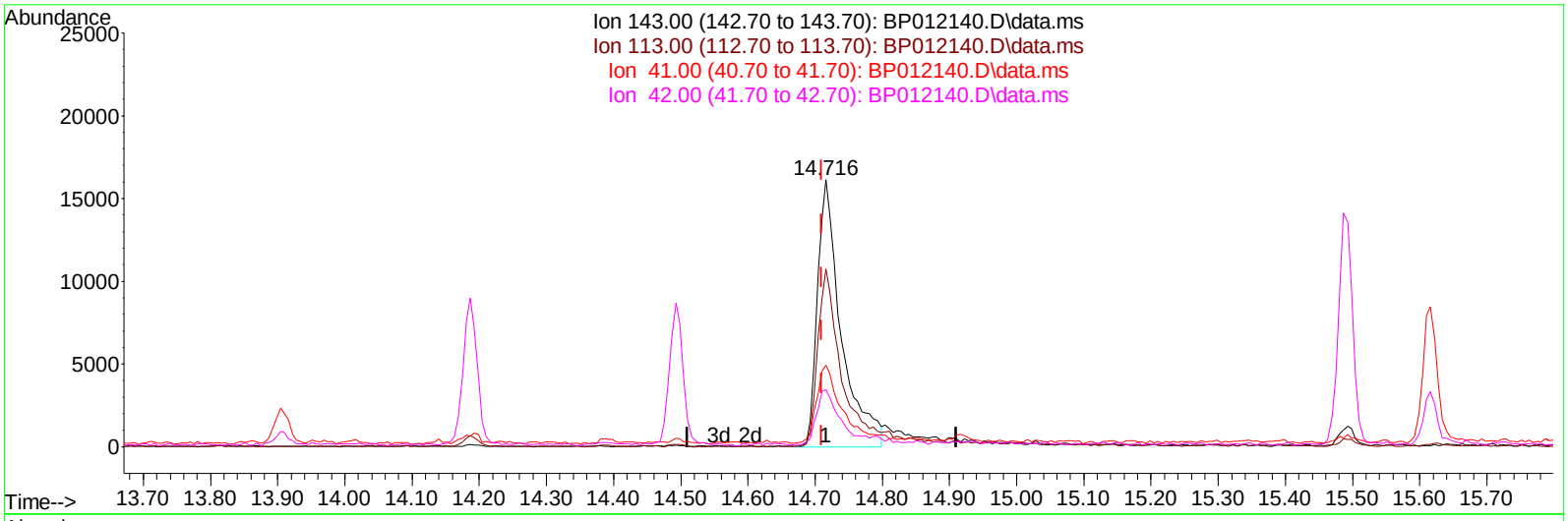

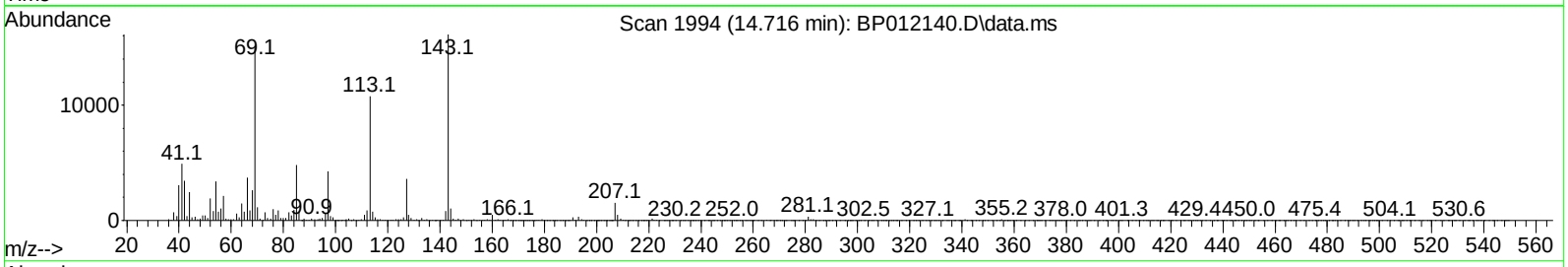

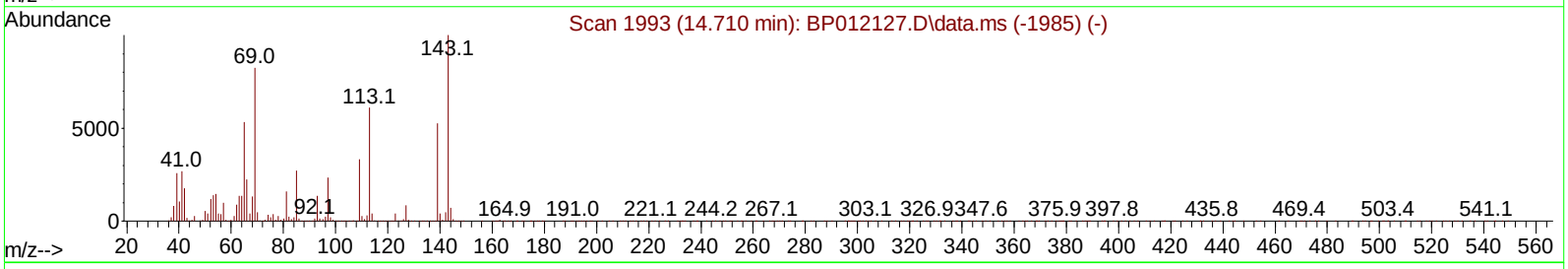

TIC: BP012140.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 5.49 ng/ul m

response 39630

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	60.80	66.67
41.00	26.60	30.46
42.00	17.60	21.41#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	255283	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	1052574	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.493	164	580984	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1204174	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1272229	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	1345707	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	28417	4.389	ng/uL	0.00
4) Pyridine-d5	3.699	84	160822	9.340	ng/ul	0.00
7) Phenol-d5	7.017	99	175741	8.560	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	417343	35.811	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	480735	29.498	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	289193	18.338	ng/ul	0.00
21) Nitrobenzene-d5	9.011	128	287808	38.716	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	286032	37.947	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.275	165	455421	31.175	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	475212	22.075	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1350415	36.701	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1628710	35.796	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	39630m	5.486	ng/ul	0.00
60) Fluorene-d10	15.493	176	1235051	37.100	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	178536	28.079	ng/ul	0.00
73) Anthracene-d10	17.357	188	2043939	39.089	ng/ul	0.00
81) Pyrene-d10	19.592	212	2479486	38.042	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	2632506	40.241	ng/ul	0.00

Target Compounds Qvalue

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

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