Quantitation Report (Qedit)

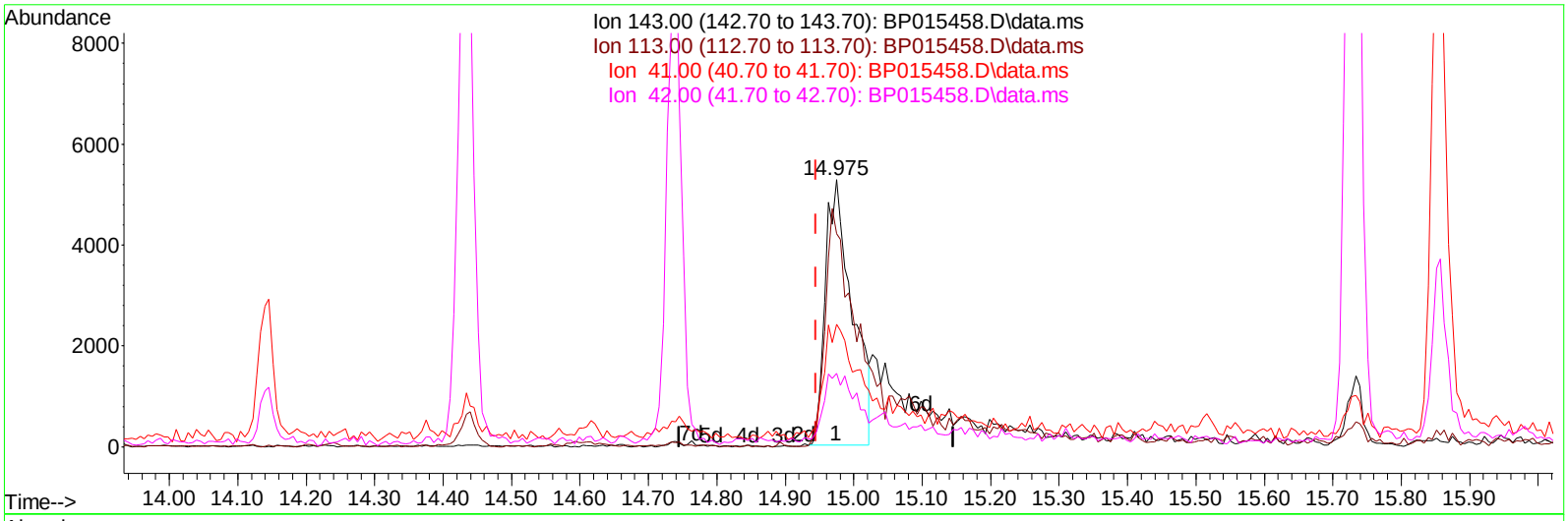
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
Data File : BP015458.D
Acq On : 09 Jun 2023 18:34
Operator : MA/JU
Sample : 03097-17
Misc :
ALS Vial : 12 Sample Multiplier: 1

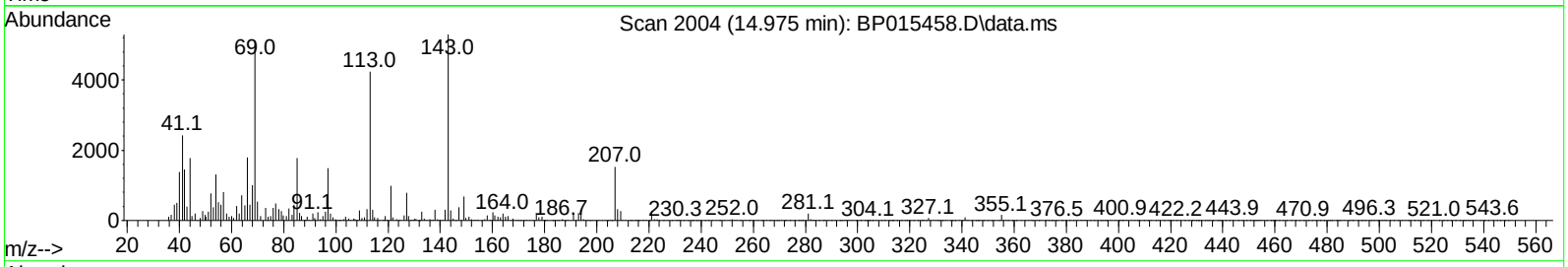
Quant Time: Jun 09 22:59:10 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 07 23:46:15 2023
Response via : Initial Calibration

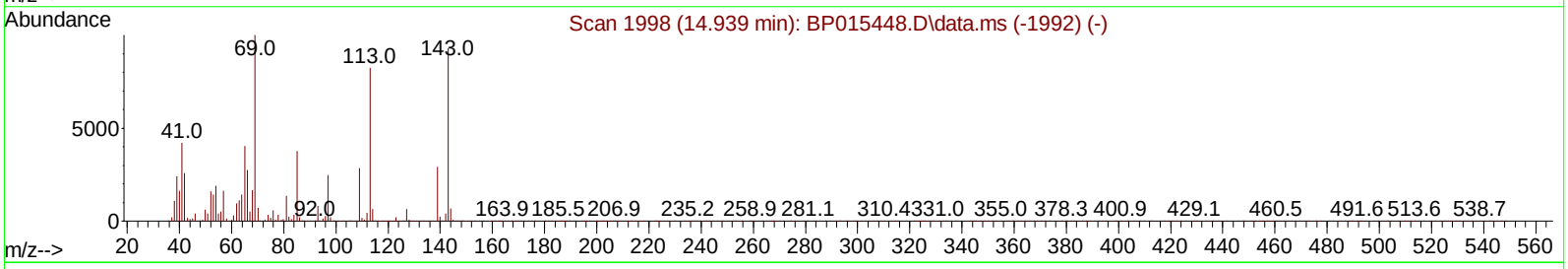
Instrument :
BNA_P
ClientSampleId :
EZEQ6

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 06/12/2023
Supervised By :mohammad ahmed 06/13/2023







TIC: BP015458.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.029) 2.65 ng/ul

response 14287

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	79.84
41.00	42.20	45.85
42.00	29.40	27.55

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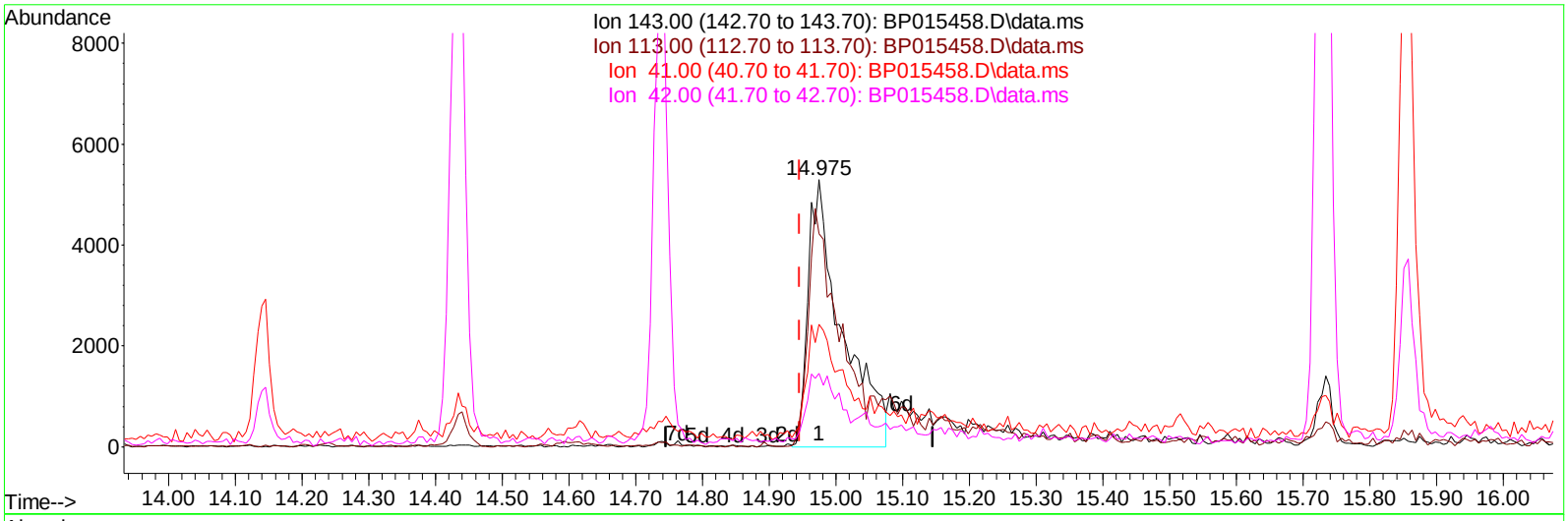

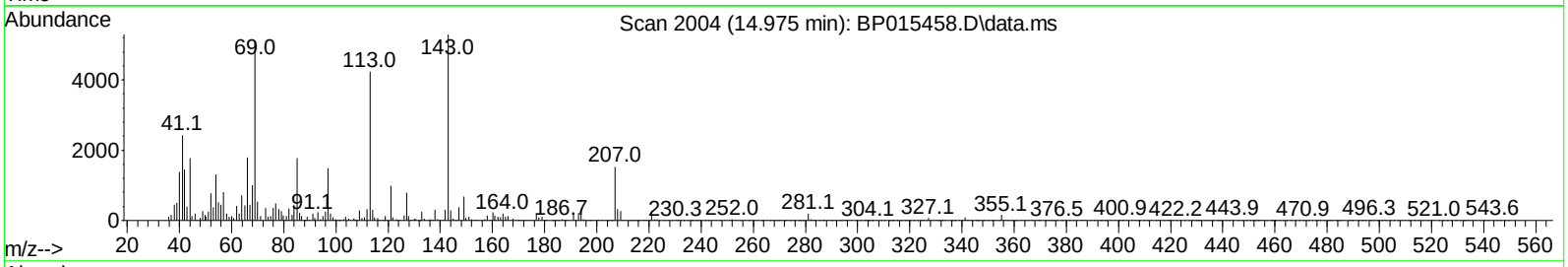

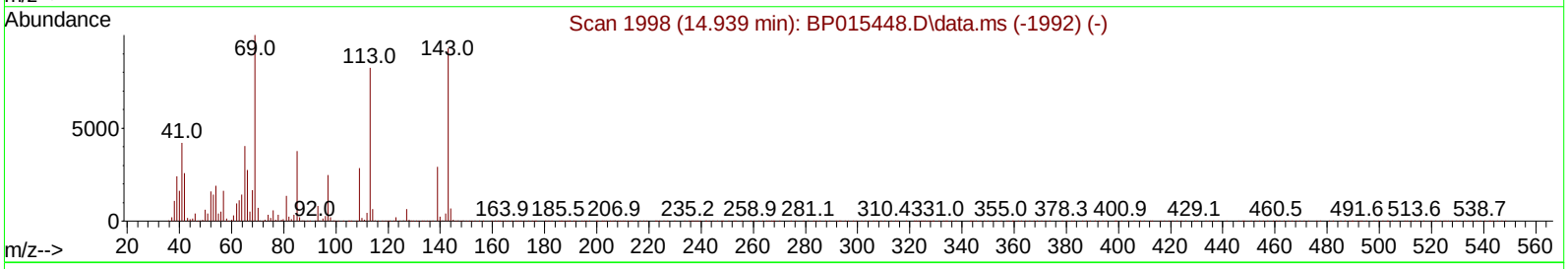

TIC: BP015458.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.975min (+ 0.029) 3.45 ng/ul m

response 18579

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	86.60	79.84
41.00	42.20	45.85
42.00	29.40	27.55

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.099	152	116948	20.000	ng/ul	0.00
20) Naphthalene-d8	10.922	136	539978	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.739	164	381396	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	902839	20.000	ng/ul	-0.01
79) Chrysene-d12	21.574	240	923534	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	977423	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	6637	2.102	ng/uL	0.00
4) Pyridine-d5	3.858	84	62716	7.645	ng/ul	0.00
7) Phenol-d5	7.252	99	66947	6.214	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	216110	33.586	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	200615	25.682	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	140724	15.915	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	140661	33.180	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	153369	31.679	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.546	165	248780	28.170	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	293448	23.149	ng/ul	0.00
46) Dimethylphthalate-d6	14.145	166	1096225	36.449	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	1163447	35.378	ng/ul	0.00
54) 4-Nitrophenol-d4	14.975	143	18579m	3.450	ng/ul	0.03
60) Fluorene-d10	15.734	176	923257	36.404	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	179676	31.414	ng/ul	0.00
73) Anthracene-d10	17.592	188	1576555	38.380	ng/ul	0.00
81) Pyrene-d10	19.816	212	2014875	40.763	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	1962759	40.021	ng/ul	-0.01
Target Compounds						Qvalue

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

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