Quantitation Report (Qedit)

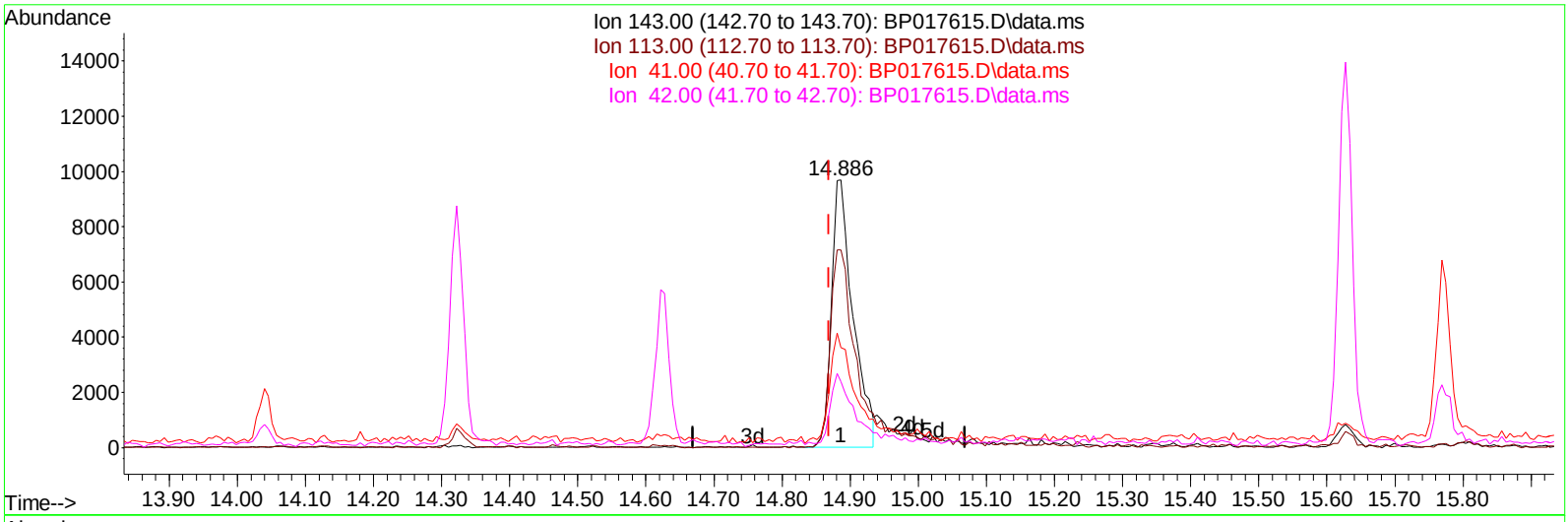
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017615.D
Acq On : 06 Oct 2023 22:46
Operator : MA/JU
Sample : 04624-19
Misc :
ALS Vial : 95 Sample Multiplier: 1

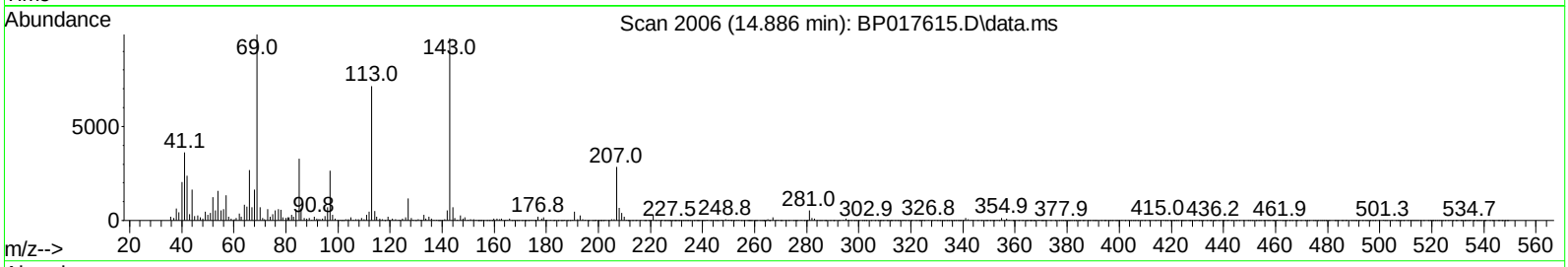
Quant Time: Oct 07 00:10:56 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 06 05:55:05 2023
Response via : Initial Calibration

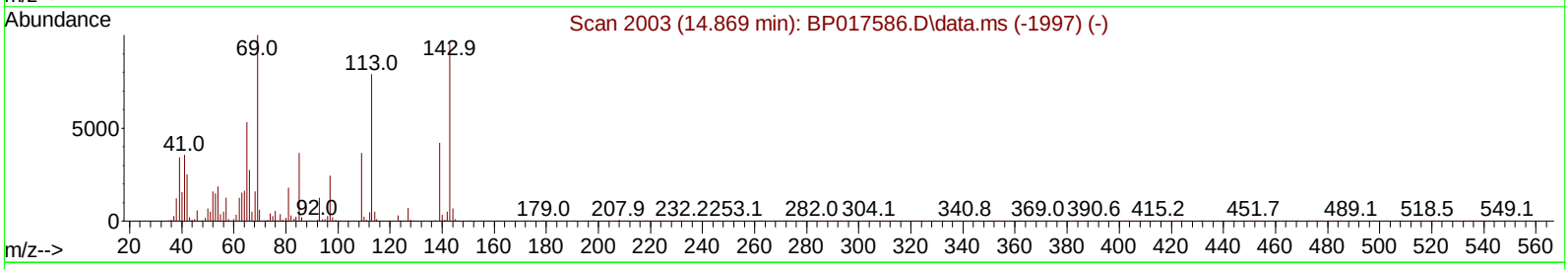
Instrument :
BNA_P
ClientSampleId :
BBHB3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/07/2023
Supervised By :mohammad ahmed 10/07/2023







TIC: BP017615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.018) 7.21 ng/ul

response 21161

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	73.80
41.00	43.80	37.31
42.00	29.00	24.73

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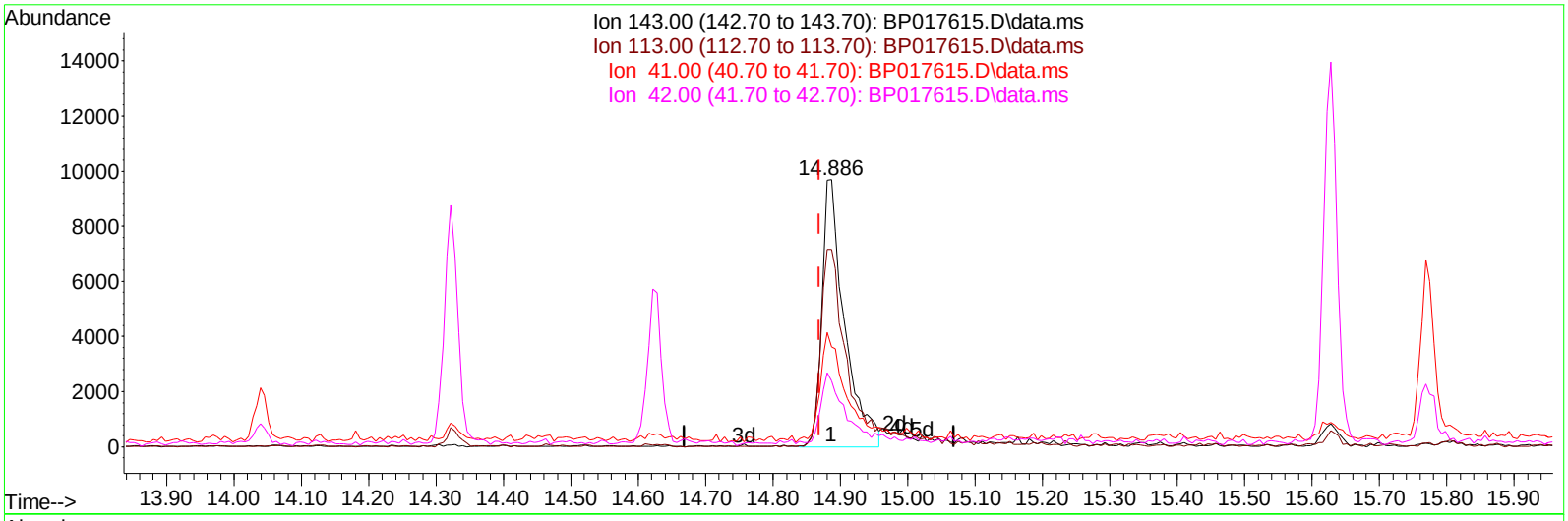

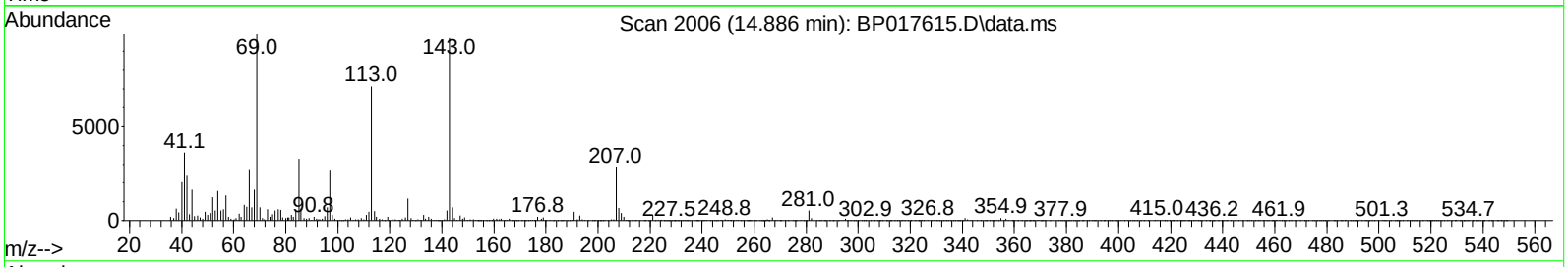

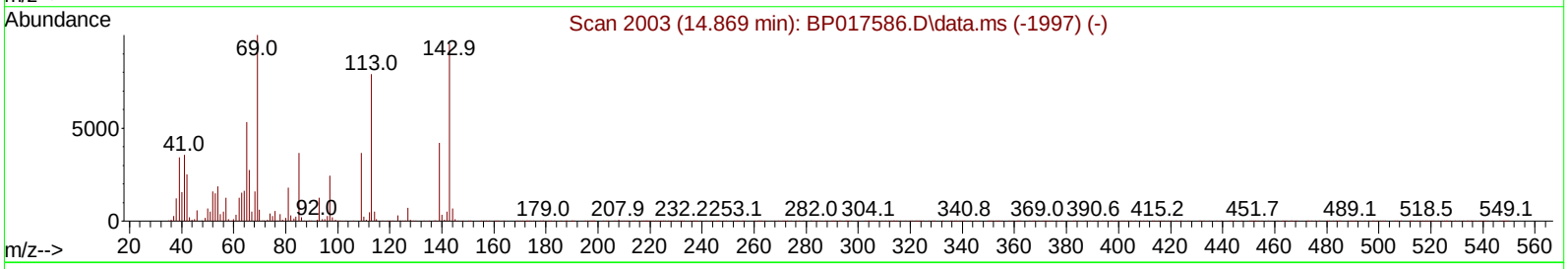

TIC: BP017615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.886min (+ 0.018) 7.68 ng/ul m

response 22526

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	81.00	73.80
41.00	43.80	37.31
42.00	29.00	24.73

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	85367	20.000	ng/ul	0.00
20) Naphthalene-d8	10.769	136	343696	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	214854	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.416	188	503389	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	529794	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	595022	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	11106	5.201	ng/uL	0.00
4) Pyridine-d5	3.728	84	86416	15.462	ng/ul	0.00
7) Phenol-d5	7.093	99	78123	10.877	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	184859	42.179	ng/ul	0.00
11) 2-Chlorophenol-d4	7.452	132	201773	35.678	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	140520	24.438	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	118207	44.705	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	125232	41.992	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	233089	42.159	ng/ul	0.00
31) 4-Chloroaniline-d4	10.940	131	292068	37.083	ng/ul	0.00
46) Dimethylphthalate-d6	14.039	166	811469	48.419	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	860987	46.590	ng/ul	0.00
54) 4-Nitrophenol-d4	14.886	143	22526m	7.675	ng/ul	0.02
60) Fluorene-d10	15.628	176	695230	48.907	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	105287	33.549	ng/ul	0.00
73) Anthracene-d10	17.516	188	1154606	49.551	ng/ul	0.00
81) Pyrene-d10	19.769	212	1469450	49.915	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	1504296	50.095	ng/ul	0.00

Target Compounds	Qvalue

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

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