Quantitation Report (Qedit)

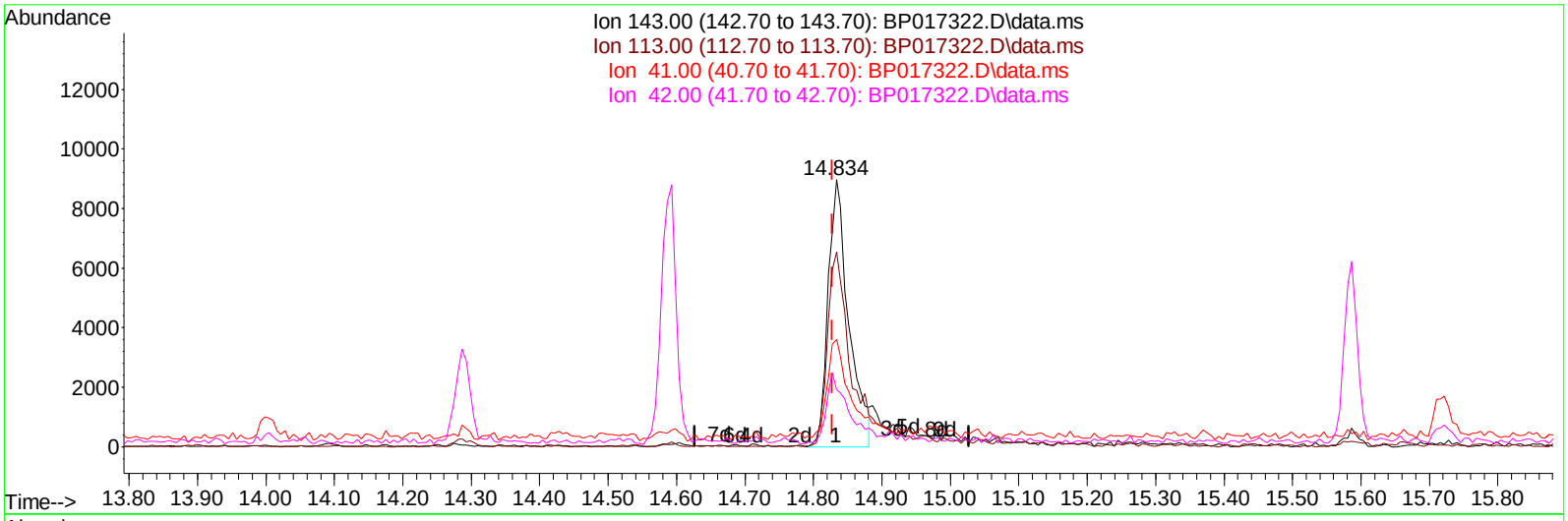
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017322.D
Acq On : 25 Sep 2023 15:44
Operator : MA/JU
Sample : 04410-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

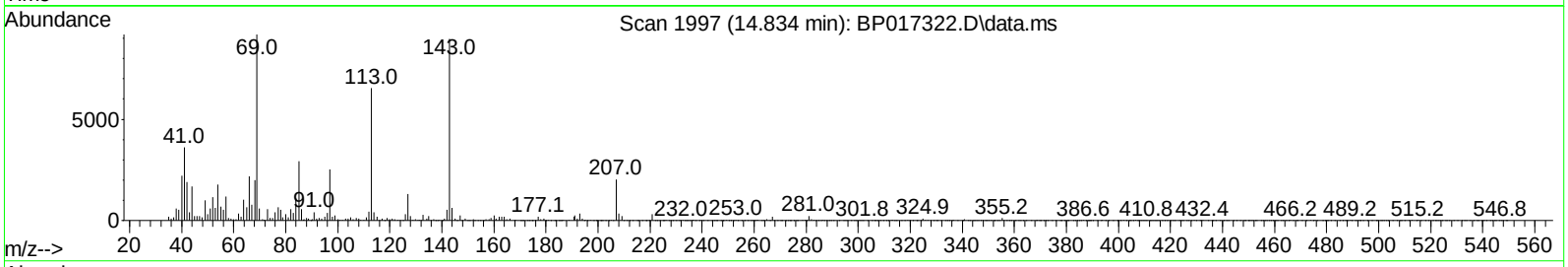
Quant Time: Sep 25 16:37:53 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 21 23:51:07 2023
Response via : Initial Calibration

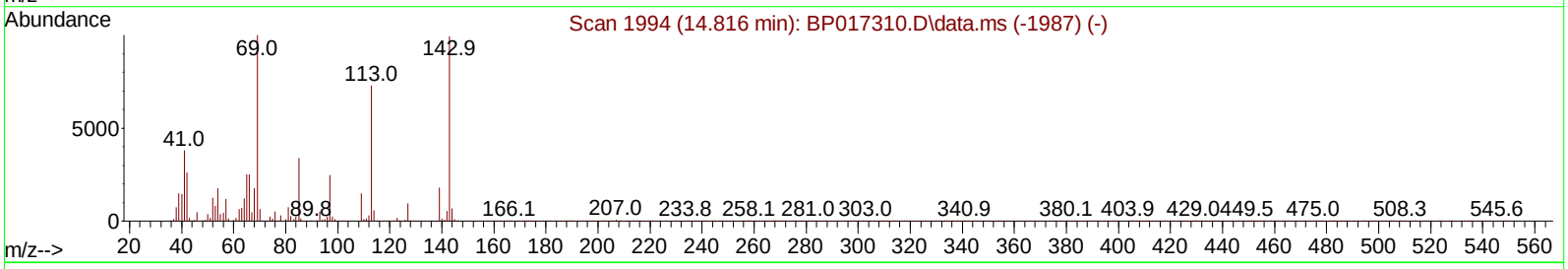
Instrument :
BNA_P
ClientSampleId :
BBJL8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/26/2023
Supervised By :mohammad ahmed 09/29/2023







TIC: BP017322.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.006) 3.41 ng/ul

response 18782

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	72.96
41.00	36.70	40.24
42.00	26.40	21.33

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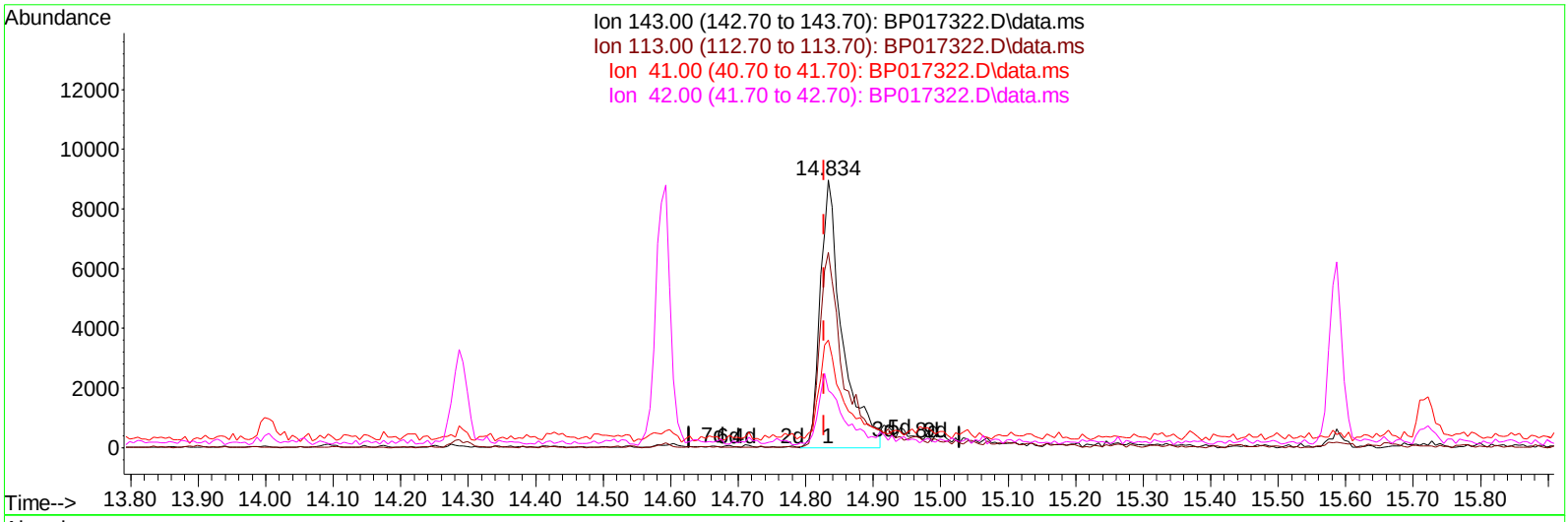

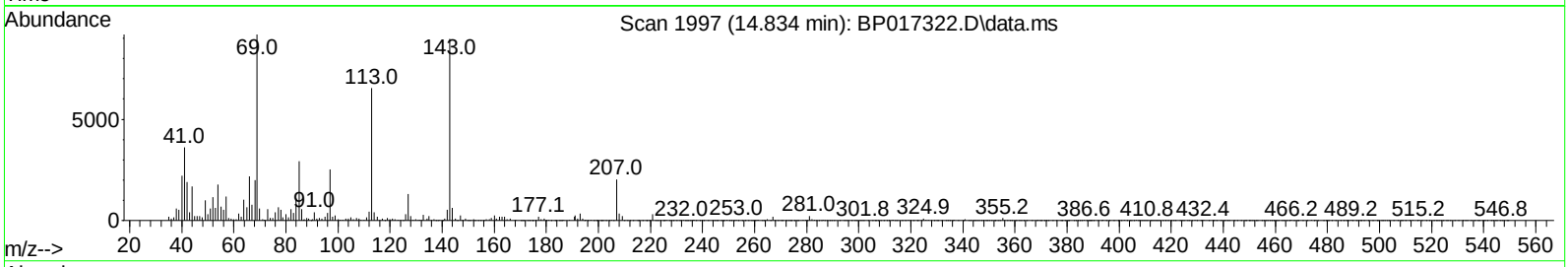

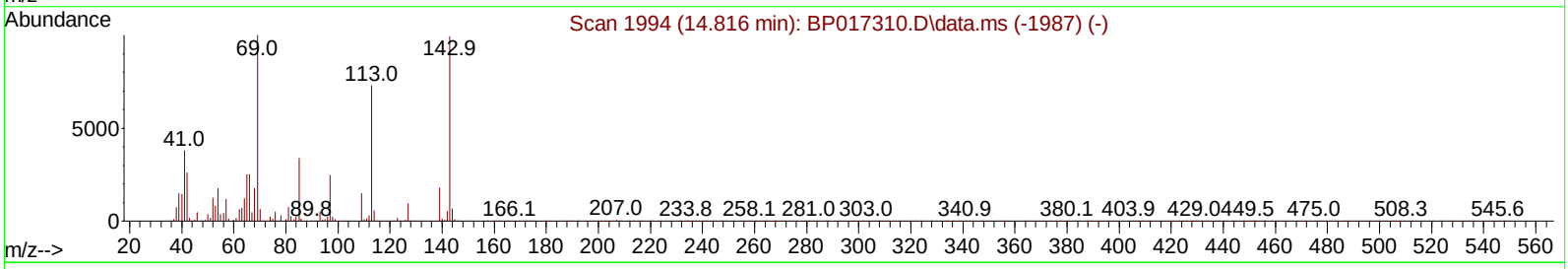

TIC: BP017322.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.834min (+ 0.006) 3.69 ng/ul m

response 20362

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.90	72.96
41.00	36.70	40.24
42.00	26.40	21.33

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
Data File : BP017322.D
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Quant Time: Sep 26 01:48:39 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Sep 21 23:51:07 2023
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	166070	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.751	136	711337	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.592	164	441717	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	991778	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	916535	20.000	ng/ul	-0.01
88) Perylene-d12	23.980	264	999331	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	13515	3.343	ng/uL	0.00
4) Pyridine-d5	3.740	84	60110	5.318	ng/ul	0.00
7) Phenol-d5	7.087	99	82875	6.076	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	91297	11.092	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.452	132	99261	9.176	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	77525	7.311	ng/ul	-0.02
21) Nitrobenzene-d5	9.128	128	52025	10.117	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.840	143	53545	9.589	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.369	165	99247	9.408	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.916	131	99207	6.492	ng/ul	-0.01
46) Dimethylphthalate-d6	14.004	166	337599	10.669	ng/ul	-0.01
49) Acenaphthylene-d8	14.287	160	375301	10.312	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.834	143	20362m	3.694	ng/ul	0.00
60) Fluorene-d10	15.587	176	317580	11.658	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.722	200	26639	4.766	ng/ul	0.00
73) Anthracene-d10	17.463	188	585878	13.320	ng/ul	0.00
81) Pyrene-d10	19.704	212	723700	14.601	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.816	264	689153	14.323	ng/ul	-0.02
Target Compounds						Qvalue

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

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