Quantitation Report (Qedit)

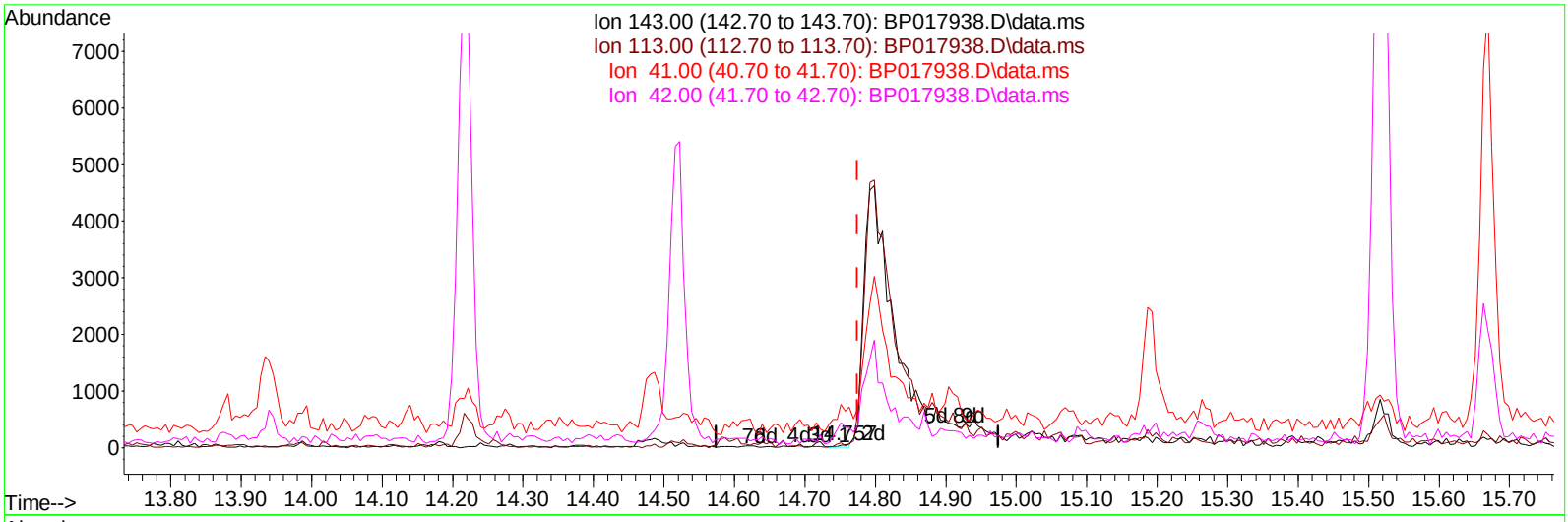
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
Data File : BP017938.D
Acq On : 07 Nov 2023 22:48
Operator : MA/JU
Sample : 05026-09
Misc :
ALS Vial : 17 Sample Multiplier: 1

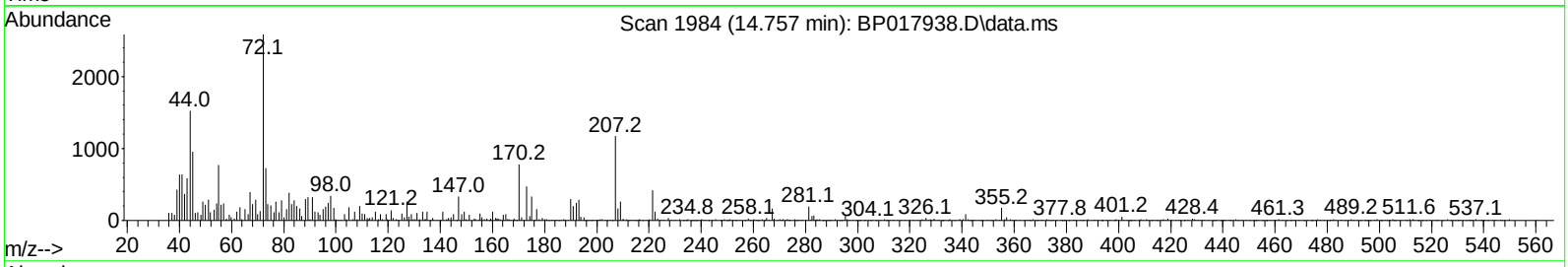
Quant Time: Nov 07 23:30:04 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 07 03:40:56 2023
Response via : Initial Calibration

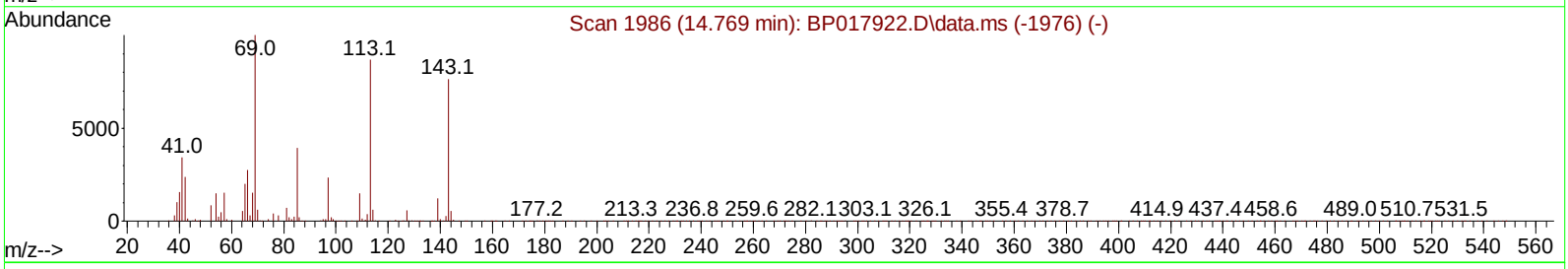
Instrument :
BNA_P
ClientSampleId :
BH336

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/09/2023
Supervised By :mohammad ahmed 11/11/2023







TIC: BP017938.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.018) 0.02 ng/ul

response 45

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	62.00#
41.00	43.00	1286.00#
42.00	28.20	736.00#

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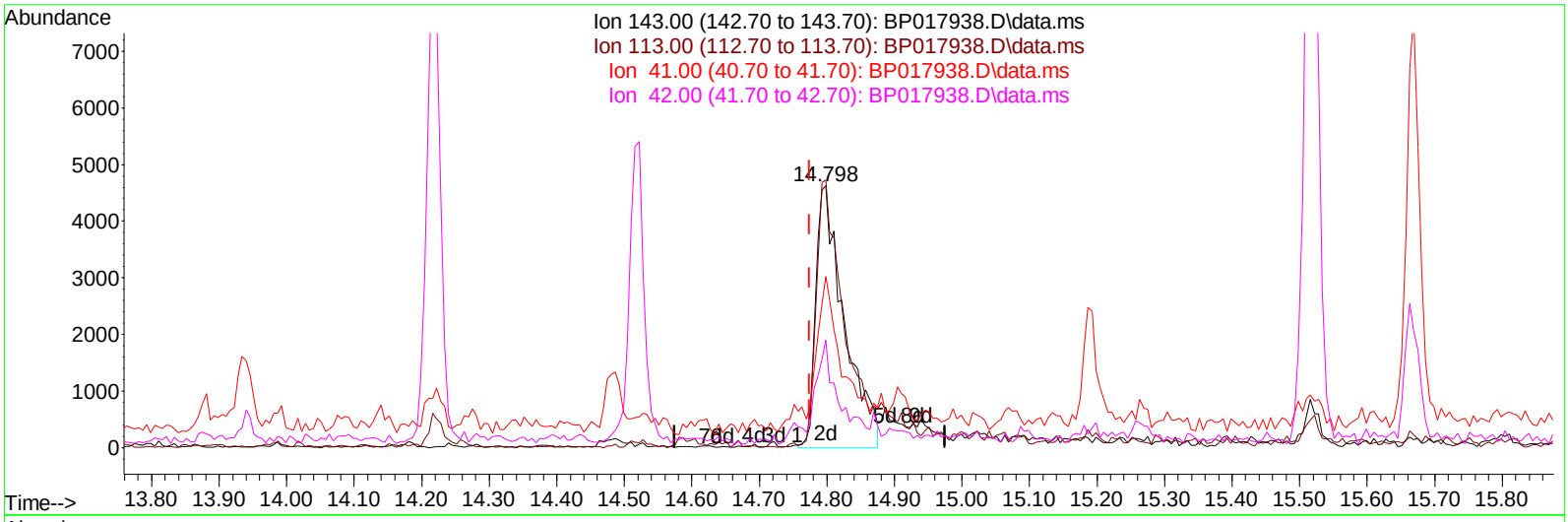

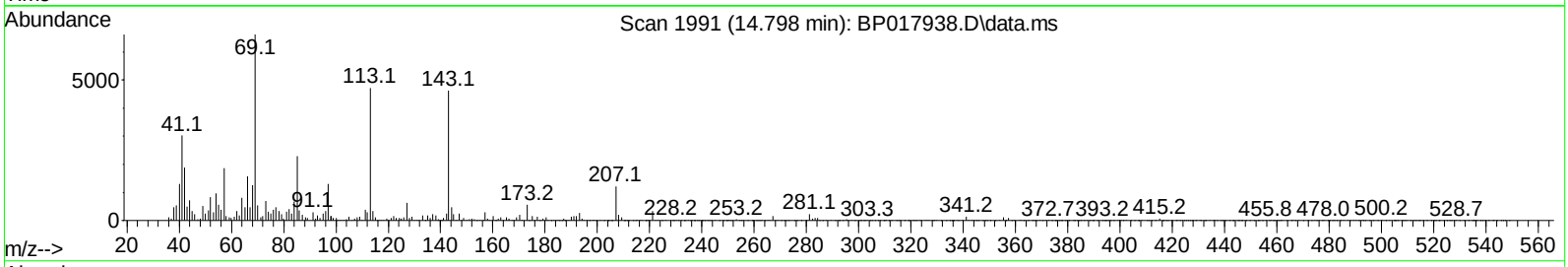

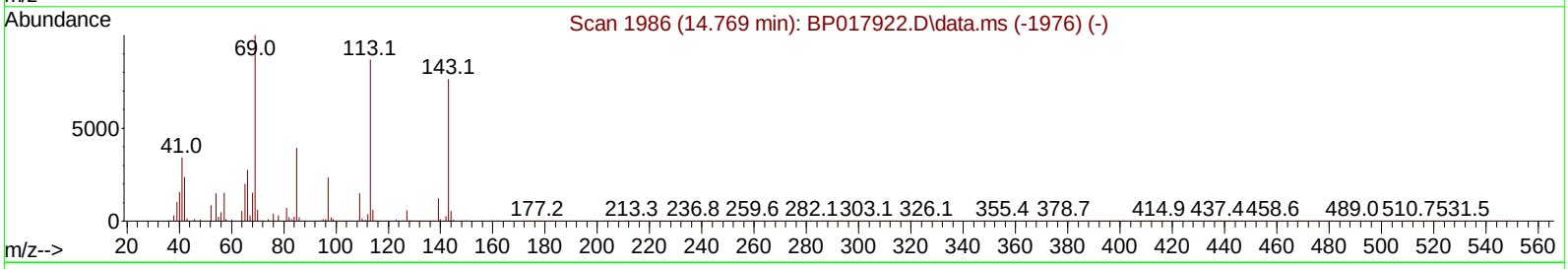

TIC: BP017938.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.798min (+ 0.024) 5.37 ng/ul m

response 13503

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	102.14
41.00	43.00	65.38#
42.00	28.20	41.02#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	69803	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	289047	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	186653	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	415738	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.410	240	393435	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	427196	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	7235	3.967	ng/uL	0.00
4) Pyridine-d5	3.687	84	47992	9.482	ng/ul	0.01
7) Phenol-d5	7.028	99	43000	6.569	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	141740	35.896	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	134006	26.642	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	88173	16.297	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	91172	39.470	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	100624	37.821	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	164282	33.682	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	203658	28.814	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	422517	25.816	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	686450	38.223	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	13503m	5.366	ng/ul	0.02
60) Fluorene-d10	15.516	176	549067	40.081	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	103374	34.002	ng/ul	0.00
73) Anthracene-d10	17.392	188	910257	42.111	ng/ul	0.00
81) Pyrene-d10	19.639	212	1132140	45.330	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1052470	44.092	ng/ul	0.00

Target Compounds Qvalue

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

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