Quantitation Report (Qedit)

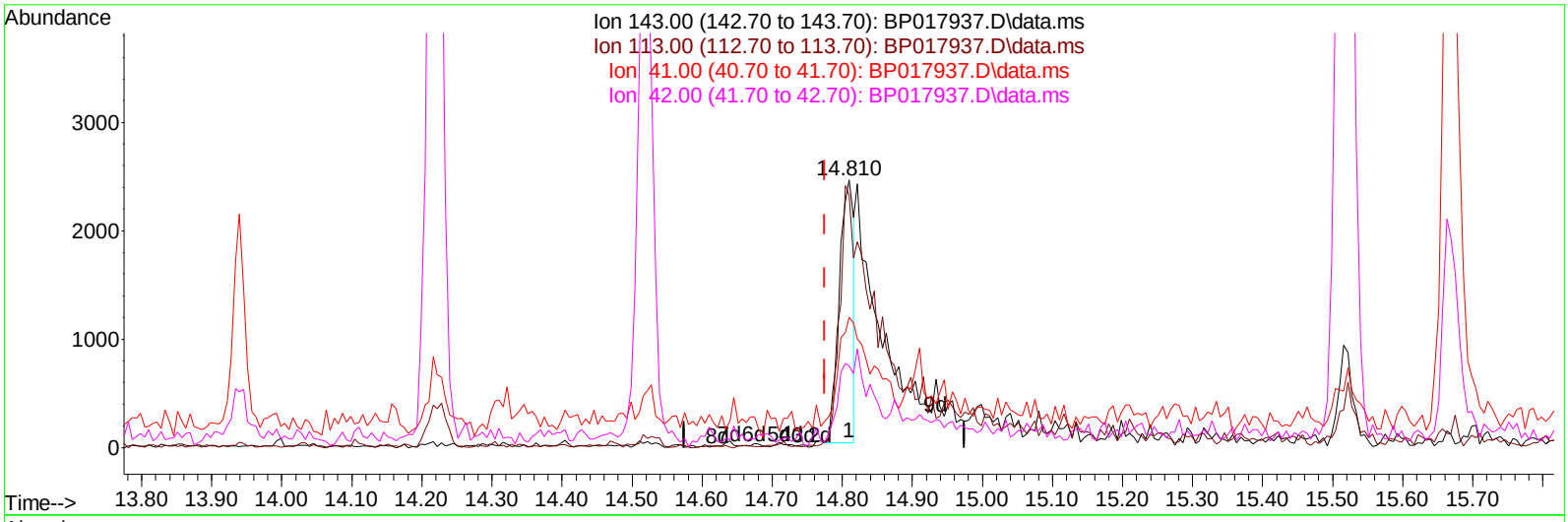
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110723\
Data File : BP017937.D
Acq On : 07 Nov 2023 22:14
Operator : MA/JU
Sample : 05026-08
Misc :
ALS Vial : 16 Sample Multiplier: 1

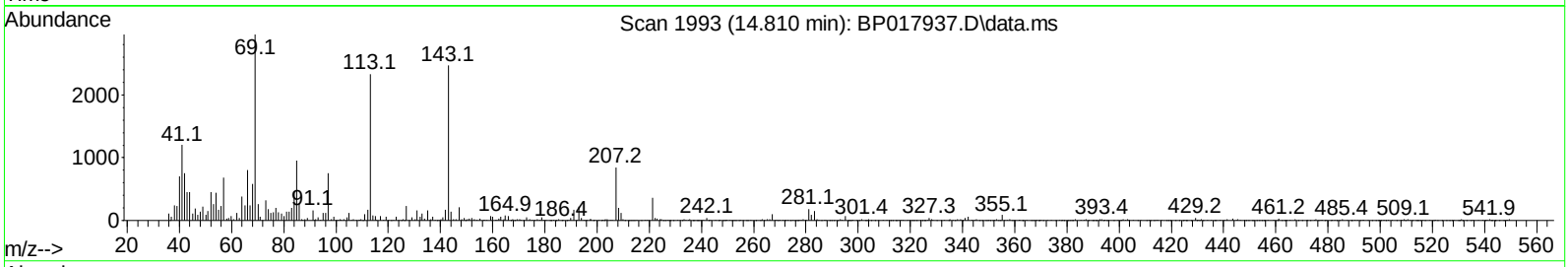
Quant Time: Nov 07 22:47:38 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110623.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 07 03:40:56 2023
Response via : Initial Calibration

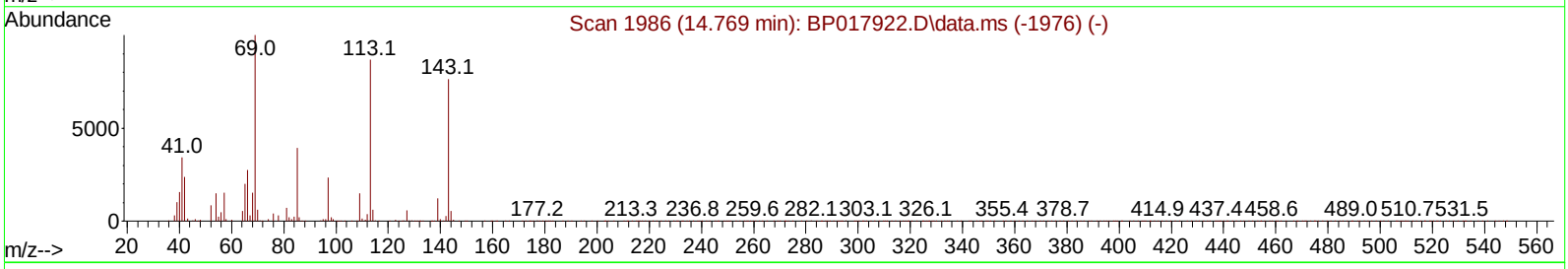
Instrument :
BNA_P
ClientSampleId :
BH338

Manual IntegrationsAPPROVED

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







TIC: BP017937.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.035) 1.47 ng/ul

response 3490

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	94.53
41.00	43.00	48.76
42.00	28.20	30.52

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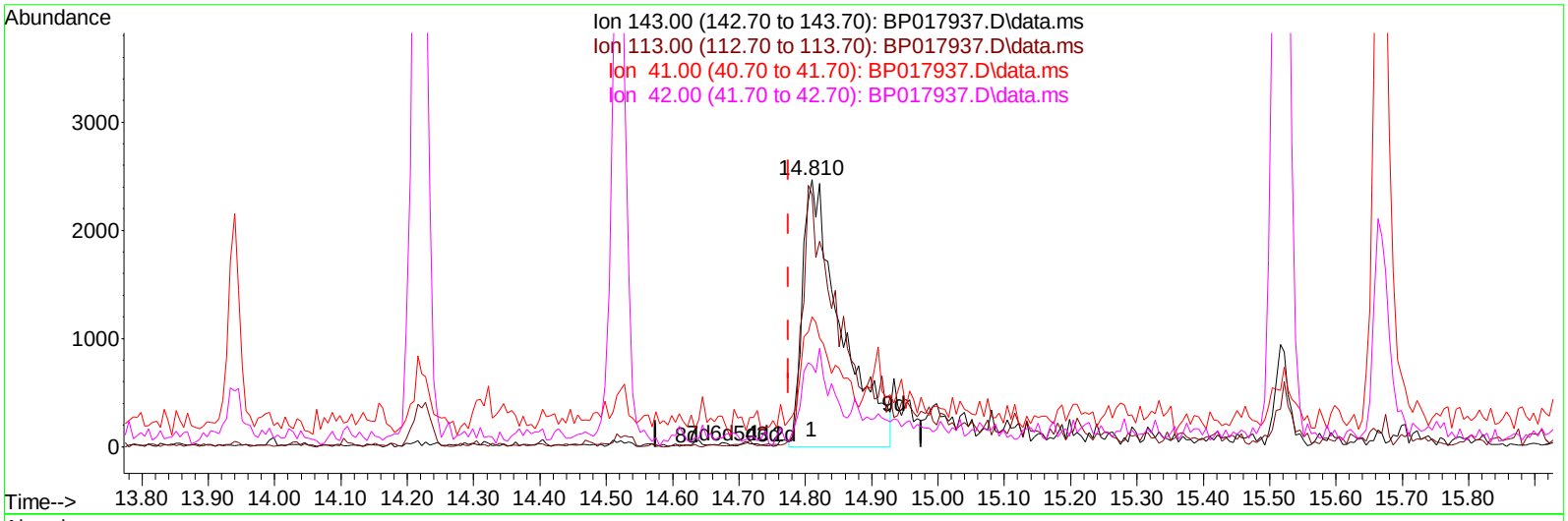

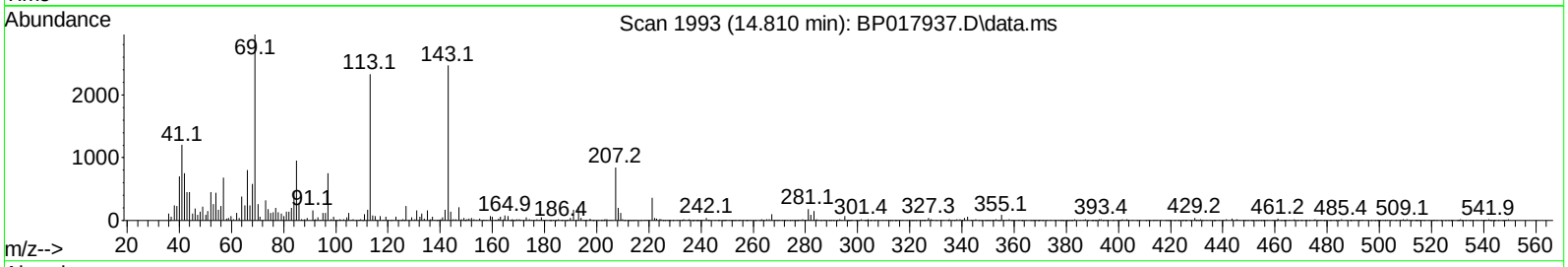

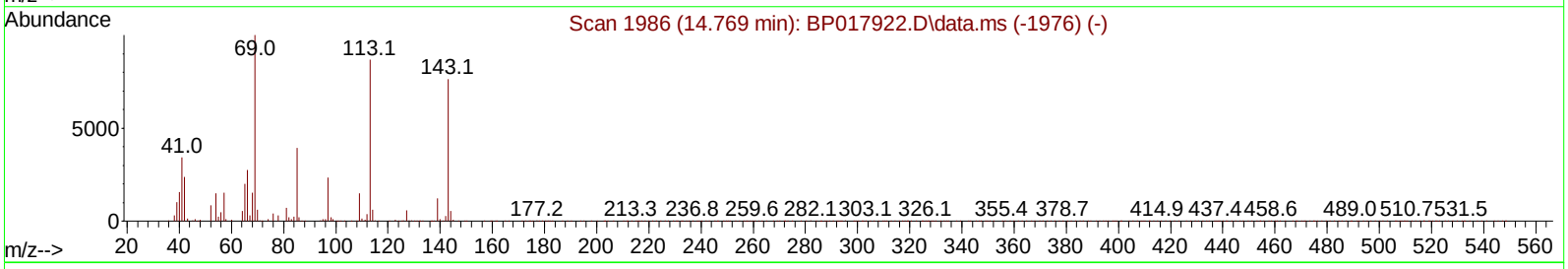

TIC: BP017937.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.035) 4.14 ng/ul m

response 9816

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	100.40	94.53
41.00	43.00	48.76
42.00	28.20	30.52

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	66323	20.000	ng/ul	0.00
20) Naphthalene-d8	10.675	136	273155	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	175809	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	393505	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	369111	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	401096	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	7863	4.537	ng/uL	0.00
4) Pyridine-d5	3.687	84	54525	11.337	ng/ul	0.01
7) Phenol-d5	7.028	99	39757	6.392	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	150716	40.172	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	136664	28.596	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	84062	16.353	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	93740	42.943	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	102076	40.599	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	160194	34.755	ng/ul	0.00
31) 4-Chloroaniline-d4	10.851	131	204224	30.576	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	644249	41.791	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	674842	39.894	ng/ul	0.00
54) 4-Nitrophenol-d4	14.810	143	9816m	4.141	ng/ul	0.04
60) Fluorene-d10	15.516	176	529181	41.013	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	97111	33.746	ng/ul	0.00
73) Anthracene-d10	17.392	188	861838	42.123	ng/ul	0.00
81) Pyrene-d10	19.645	212	1046301	44.654	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	986305	44.009	ng/ul	0.00

Target Compounds Qvalue

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

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