

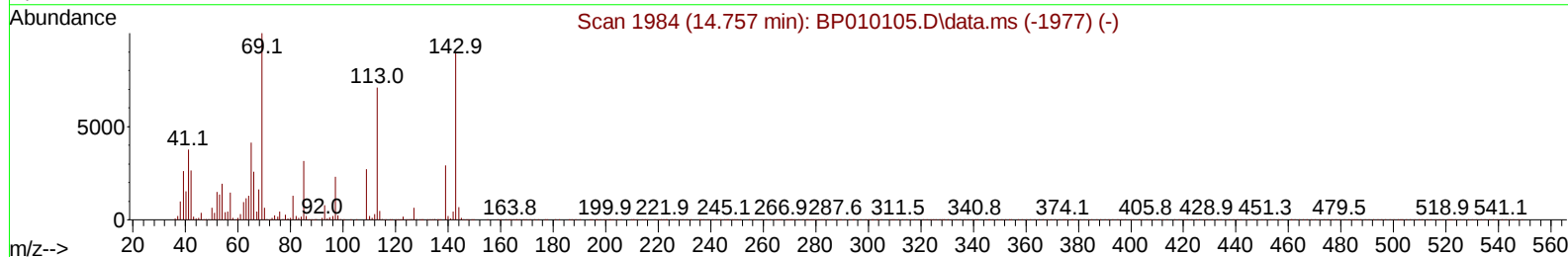
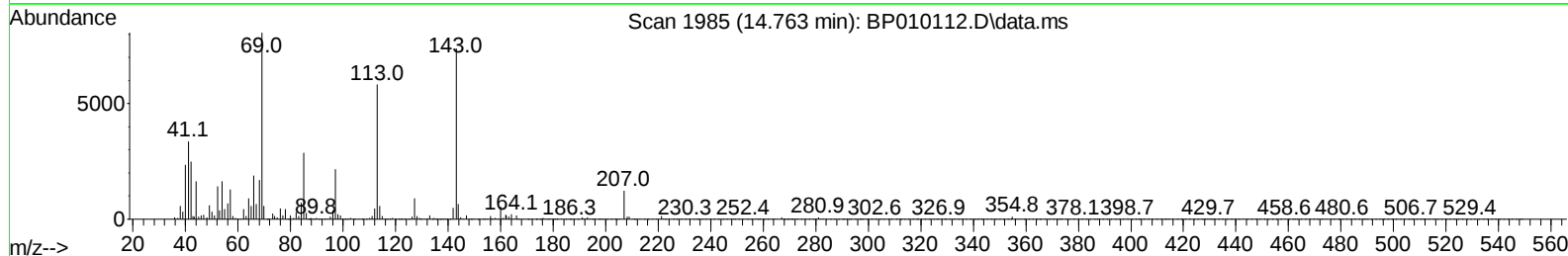
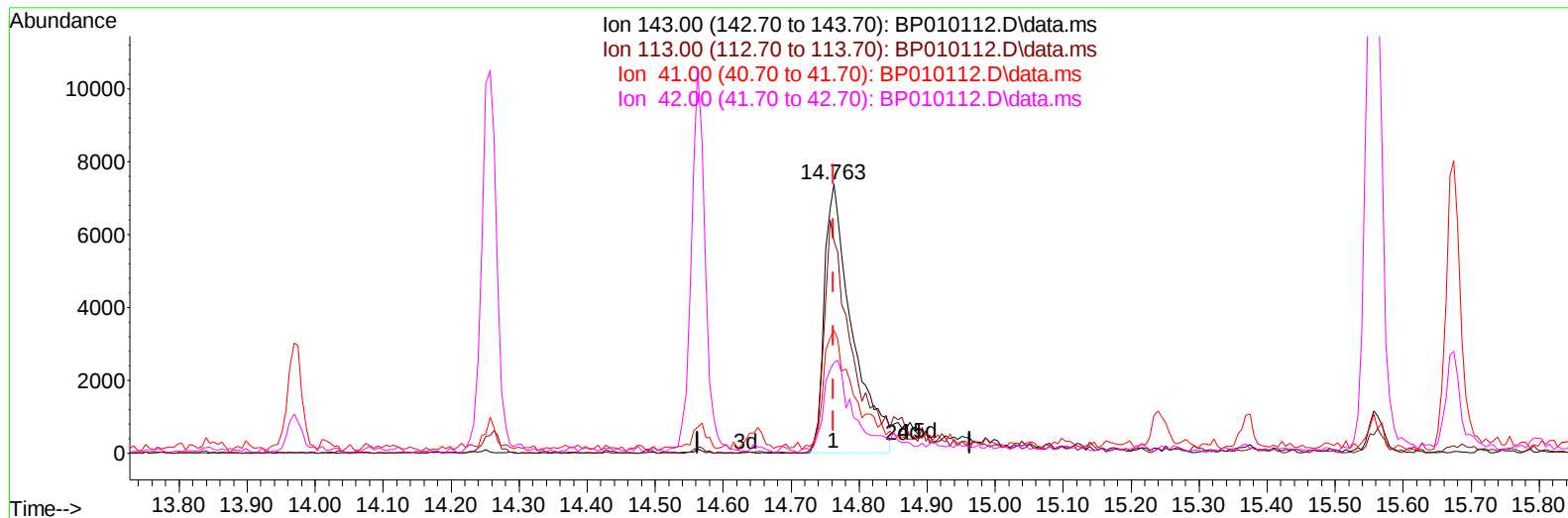
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010112.D
Acq On : 26 Apr 2022 12:58
Operator : CG/JU
Sample : N2517-03
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFH8

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/27/2022
Supervised By :Yogesh Patel 04/28/2022

Quant Time: Apr 27 00:45:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration



TIC: BP010112.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 3.33 ng/u1

response 20980

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	78.76#
41.00	23.20	45.60#
42.00	18.00	33.55#

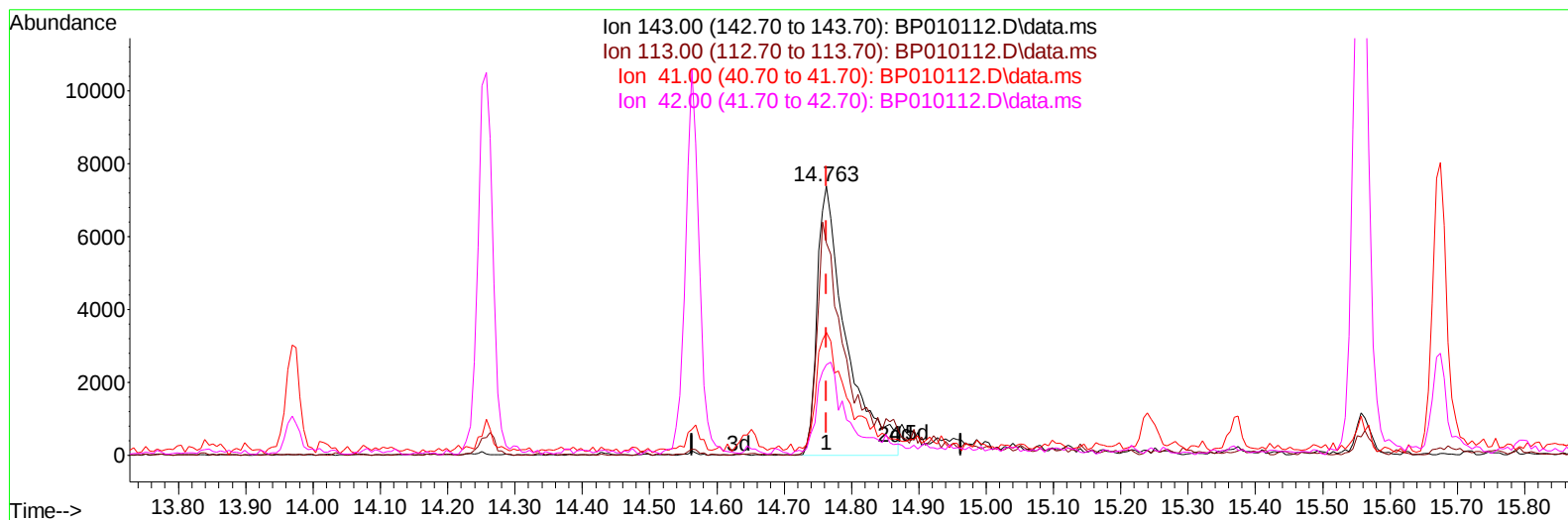
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	144231	20.000	ng/ul	0.00
20) Naphthalene-d8	10.739	136	623736	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	439819	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.316	188	921168	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	799530	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264	627225	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	16209	4.901	ng/uL	0.00
4) Pyridine-d5	3.793	84	37825	4.065	ng/ul	0.00
7) Phenol-d5	7.093	99	78124	5.955	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.257	67	243533	30.936	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	224898	23.147	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	156183	14.369	ng/ul	-0.02
21) Nitrobenzene-d5	9.092	128	154209	31.085	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	151968	28.049	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	259921	24.643	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	333911	22.079	ng/ul	-0.01
46) Dimethylphthalate-d6	13.969	166	1110107	32.447	ng/ul	-0.01
49) Acenaphthylene-d8	14.257	160	1225617	29.599	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	22158m	3.518	ng/ul	0.00
60) Fluorene-d10	15.557	176	960573	32.639	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.674	200	131579	22.590	ng/ul	0.00
73) Anthracene-d10	17.416	188	1687685	37.870	ng/ul	0.00
81) Pyrene-d10	19.645	212	1944716	42.770	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1278040	38.867	ng/ul	0.00

Target Compounds Qvalue

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

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