

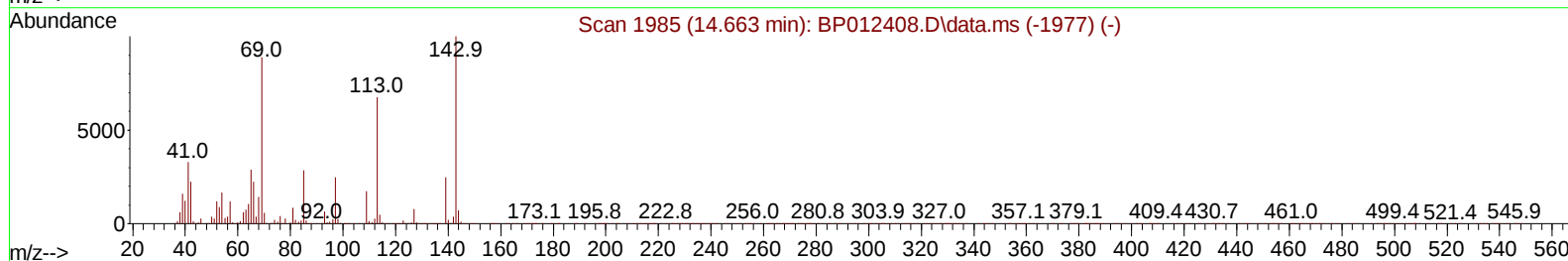
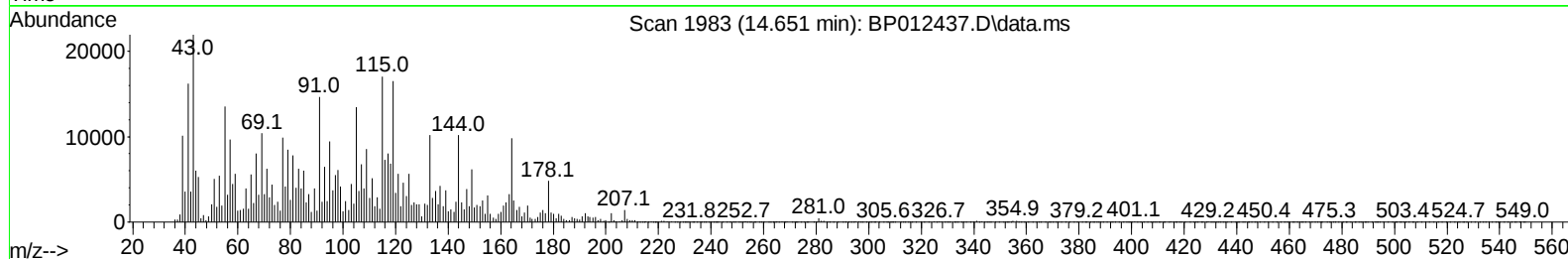
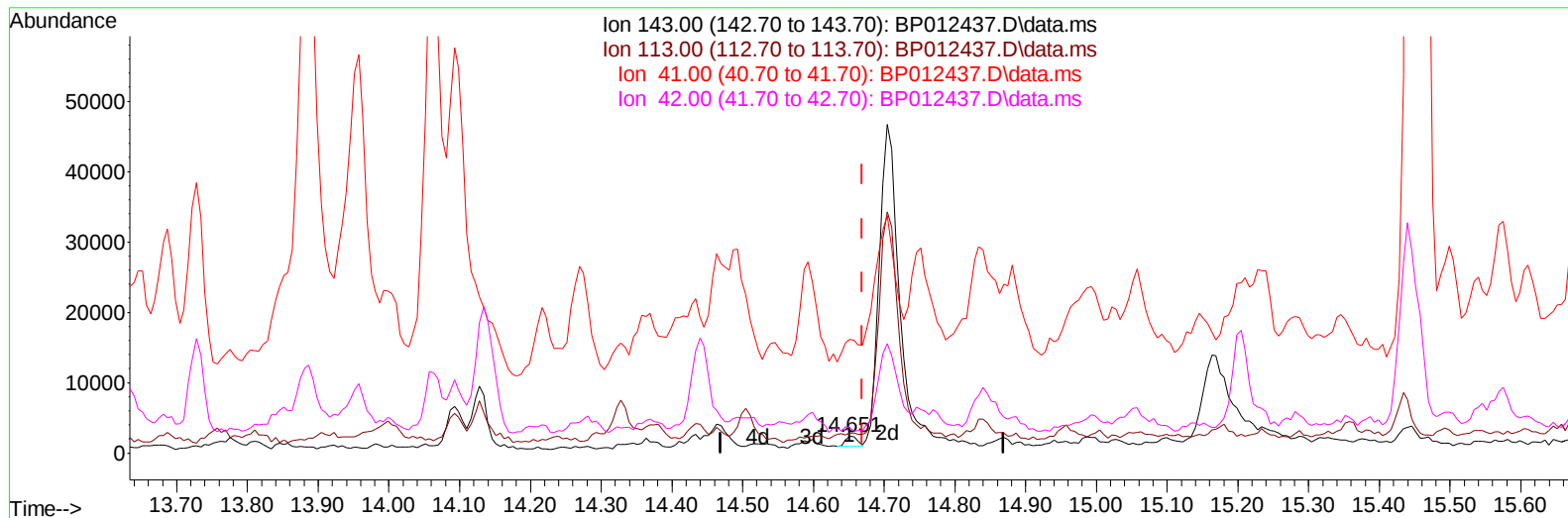
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012437.D
Acq On : 01 Nov 2022 20:45
Operator : CG/JU
Sample : N5300-02
Misc :
ALS Vial : 58 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA73

Manual IntegrationsAPPROVED

Quant Time: Nov 01 23:01:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 01 01:03:58 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/02/2022
Supervised By :mohammad ahmed 11/02/2022



TIC: BP012437.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.651min (-0.018) 0.15 ng/u1

response 1636

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	121.40#
41.00	34.50	685.36#
42.00	23.60	150.68#

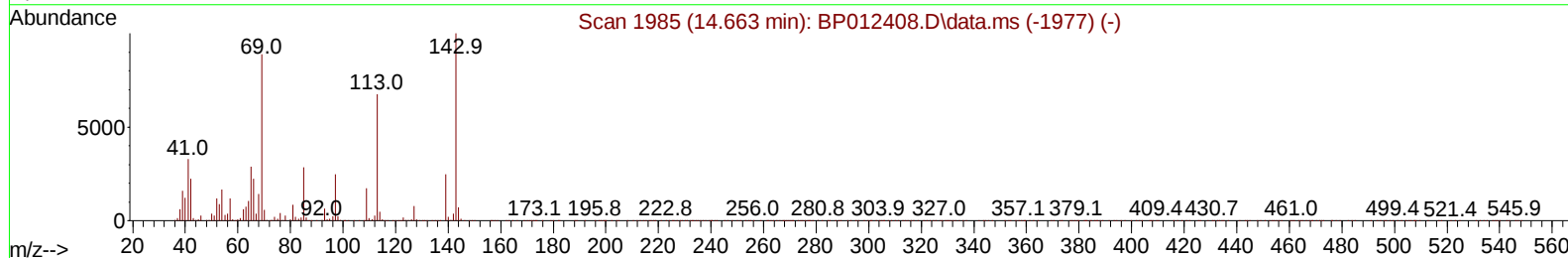
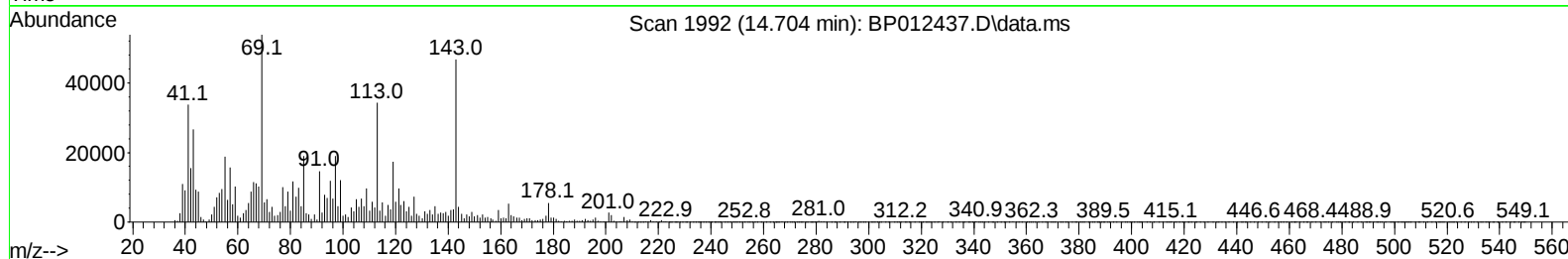
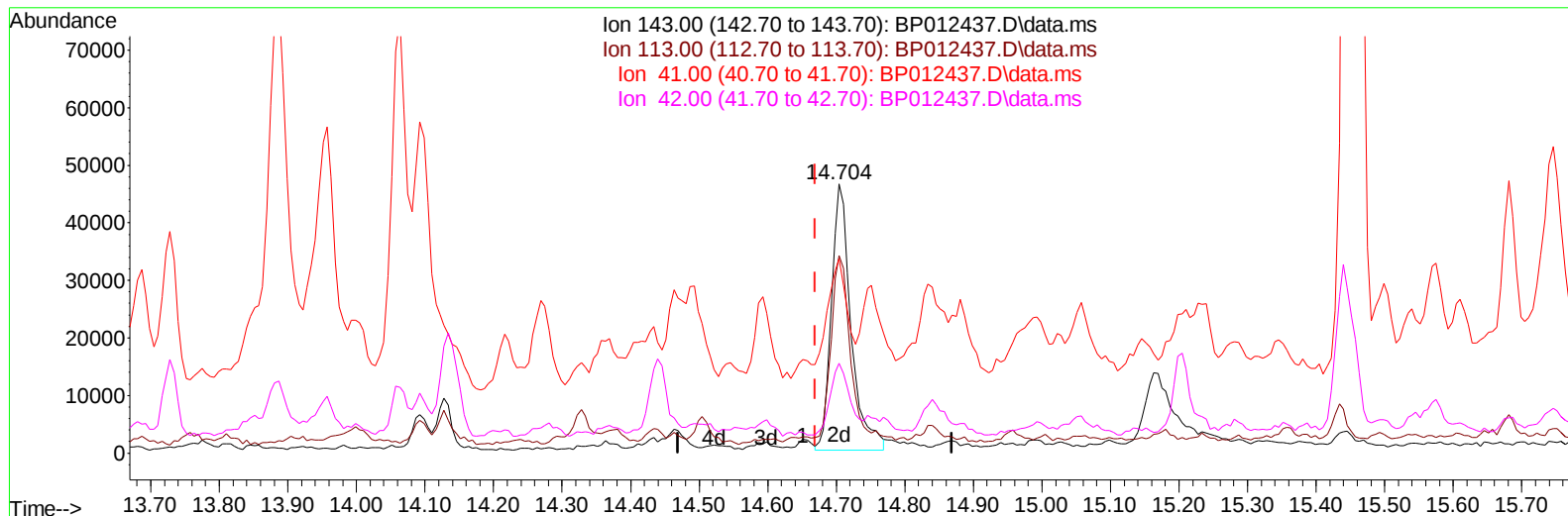
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TIC: BP012437.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.035) 8.66 ng/ul m

response 92123

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.10	73.39
41.00	34.50	72.20#
42.00	23.60	33.31#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	295439	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	1290473	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	820621	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1736245	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1600177	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1346755	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	28316	3.859	ng/uL	0.00
4) Pyridine-d5	3.664	84	137588	6.528	ng/ul	0.00
7) Phenol-d5	6.969	99	226883	8.562	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	536577	33.920	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	580652	30.006	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	394646	18.830	ng/ul	0.00
21) Nitrobenzene-d5	8.958	128	362488	43.316	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	390223	44.926	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	673200	35.763	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	589060	21.017	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2144175	38.169	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2424537	37.767	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	92123m	8.665	ng/ul	0.04
60) Fluorene-d10	15.440	176	1738346	36.143	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	266425	29.473	ng/ul	0.00
73) Anthracene-d10	17.304	188	2894257	38.688	ng/ul	0.00
81) Pyrene-d10	19.545	212	3266763	40.662	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	2528669	38.427	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.270	88	230872	29.597	ng/uL	100
8) Phenol	6.999	94	132992	4.827	ng/ul#	92
26) 2,4-Dimethylphenol	9.781	107	177030	7.741	ng/ul	97
30) Naphthalene	10.658	128	17536619	256.410	ng/ul	97
36) 2-Methylnaphthalene	12.257	142	2712353	57.847	ng/ul	100
37) 1-Methylnaphthalene	12.481	142	3290665	70.273	ng/ul	99
43) 1,1'-Biphenyl	13.269	154	104829	1.737	ng/ul	96
52) Acenaphthene	14.504	153	926164	18.305	ng/ul	99

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

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