

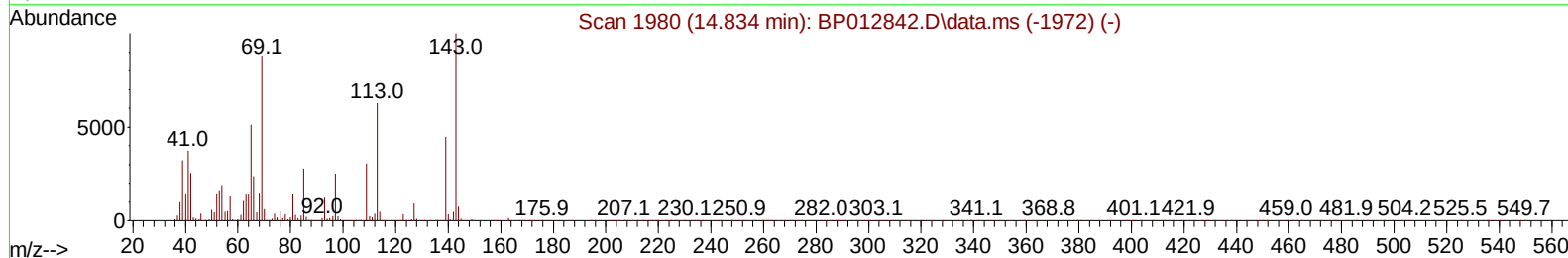
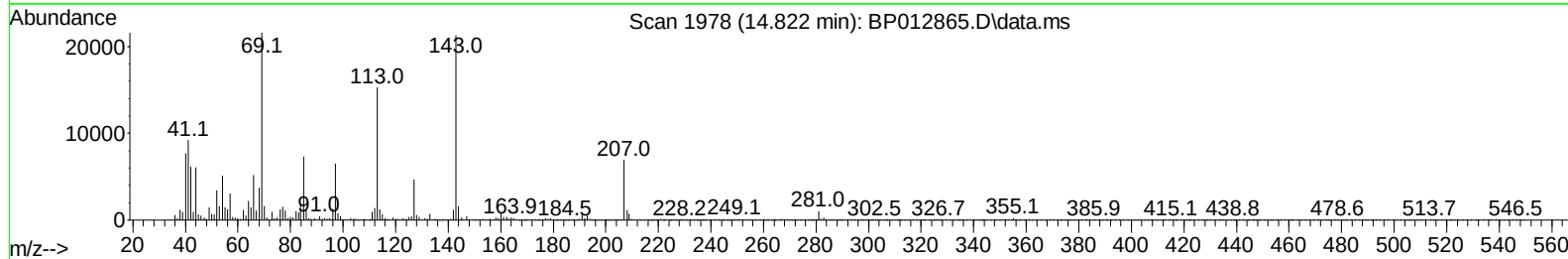
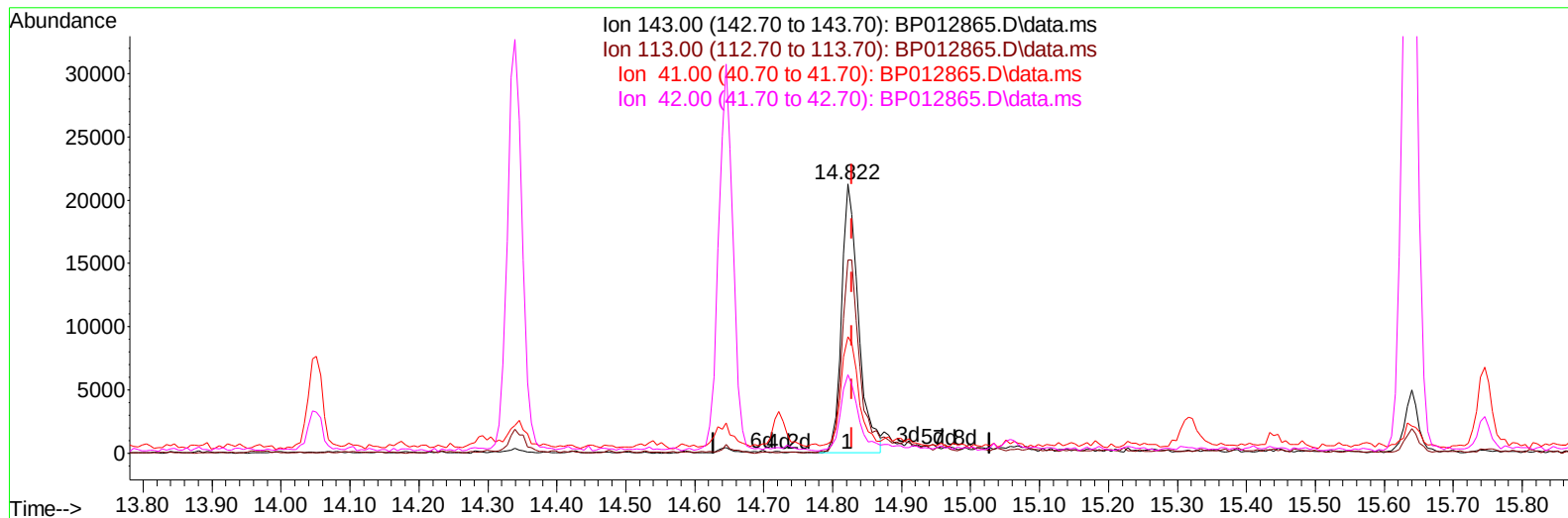
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012865.D
 Acq On : 29 Nov 2022 22:36
 Operator : CG/JU
 Sample : N5725-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB56

Manual IntegrationsAPPROVED

Quant Time: Nov 29 23:26:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/30/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012865.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (-0.006) 1.64 ng/ul

response 36108

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	71.77
41.00	37.00	43.25
42.00	24.70	29.05

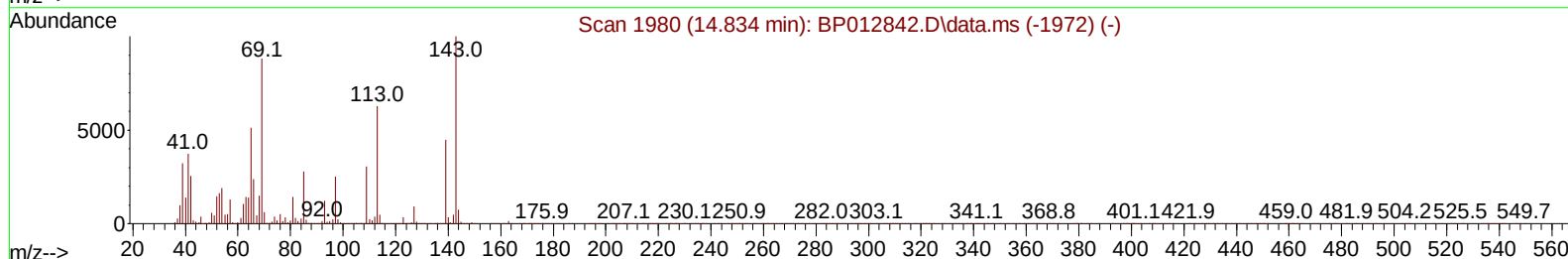
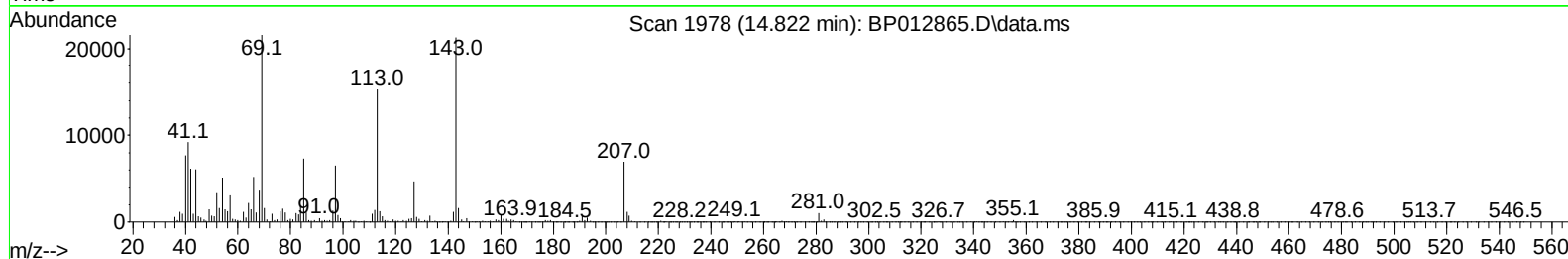
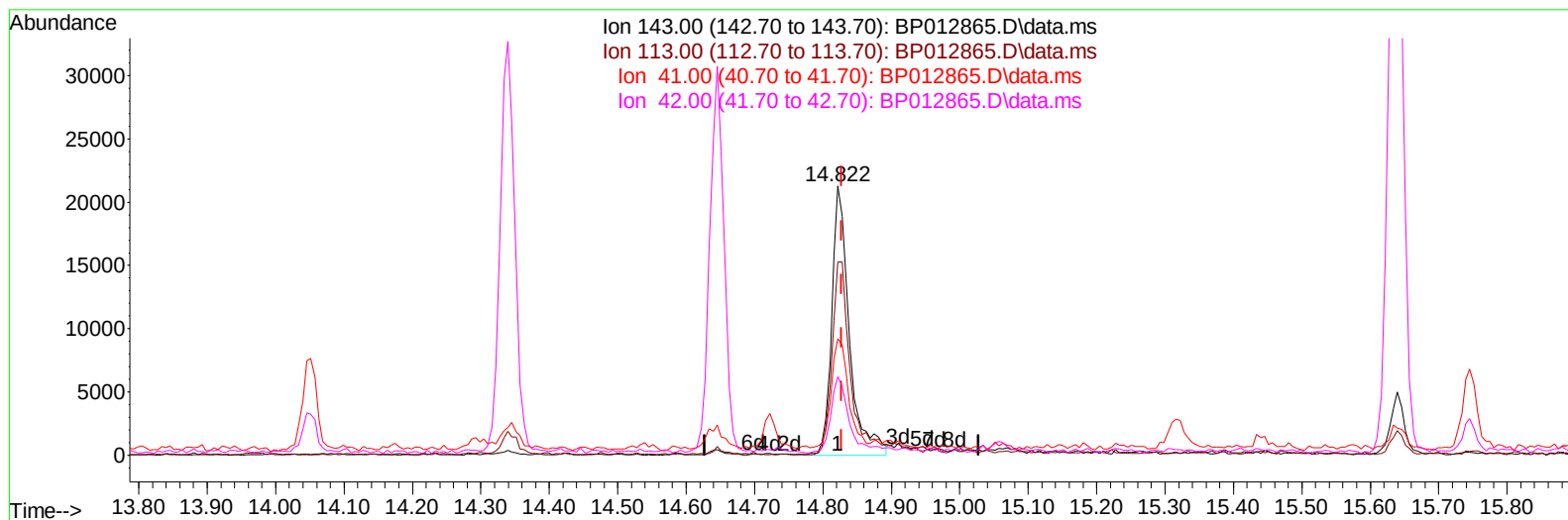
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TIC: BP012865.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.822min (-0.006) 1.73 ng/ul m

response 37908

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	71.77
41.00	37.00	43.25
42.00	24.70	29.05

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	577054	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2615522	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1789242	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	3702790	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	2505106	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	2065389	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	55863	4.148	ng/uL	0.00
4) Pyridine-d5	3.828	84	141246	3.505	ng/ul	0.01
7) Phenol-d5	7.158	99	239359	5.048	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	904209	30.946	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	517207	14.314	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	382243	10.005	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128	588103	36.737	ng/ul	0.00
24) 2-Nitrophenol-d4	9.905	143	272040	16.004	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	602803	15.199	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	522932	9.628	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	3753476	30.477	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	4778018	33.250	ng/ul	0.00
54) 4-Nitrophenol-d4	14.822	143	37908m	1.726	ng/ul	0.00
60) Fluorene-d10	15.640	176	3912637	36.276	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	99562	5.626	ng/ul	0.00
73) Anthracene-d10	17.498	188	6313490	38.735	ng/ul	0.00
81) Pyrene-d10	19.728	212	6519289	44.787	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	3962988	39.063	ng/ul	0.00

Target Compounds Qvalue

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

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