

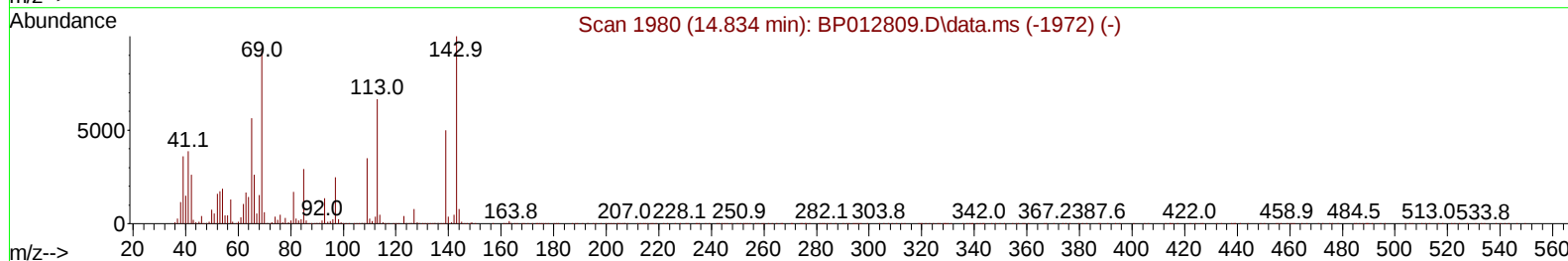
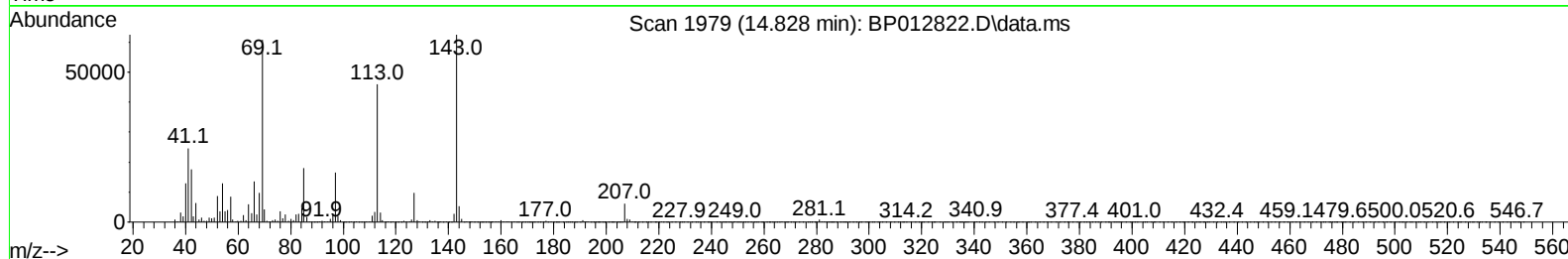
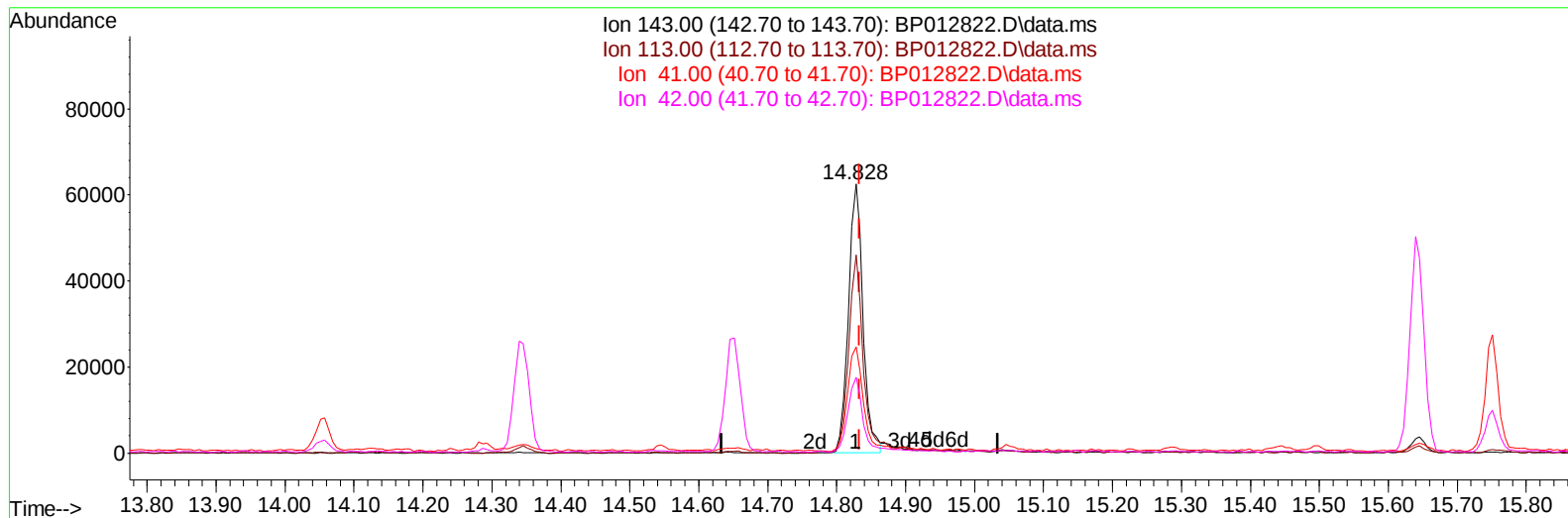
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
Data File : BP012822.D
Acq On : 28 Nov 2022 18:13
Operator : CG/JU
Sample : N5666-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHB32

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/29/2022
Supervised By :mohammad ahmed 11/30/2022

Quant Time: Nov 28 22:42:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Nov 27 22:55:08 2022
Response via : Initial Calibration



TIC: BP012822.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.828min (-0.006) 4.82 ng/u1****response 91835**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	73.64
41.00	37.00	39.44
42.00	24.70	28.22

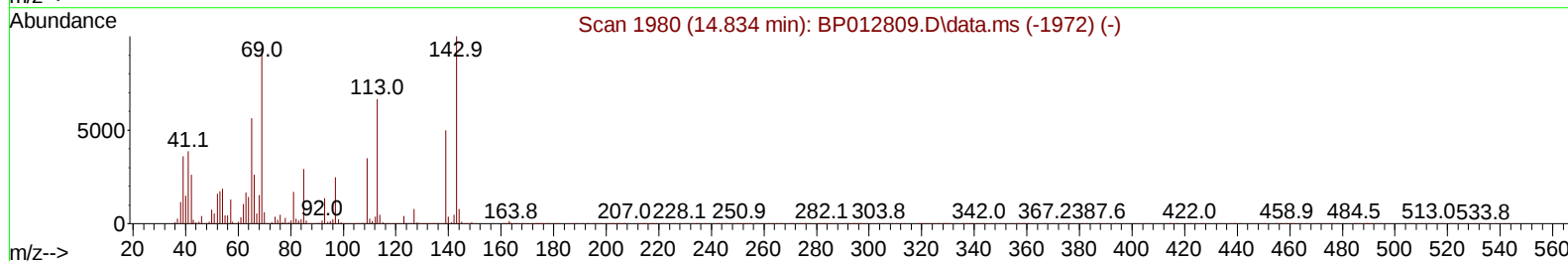
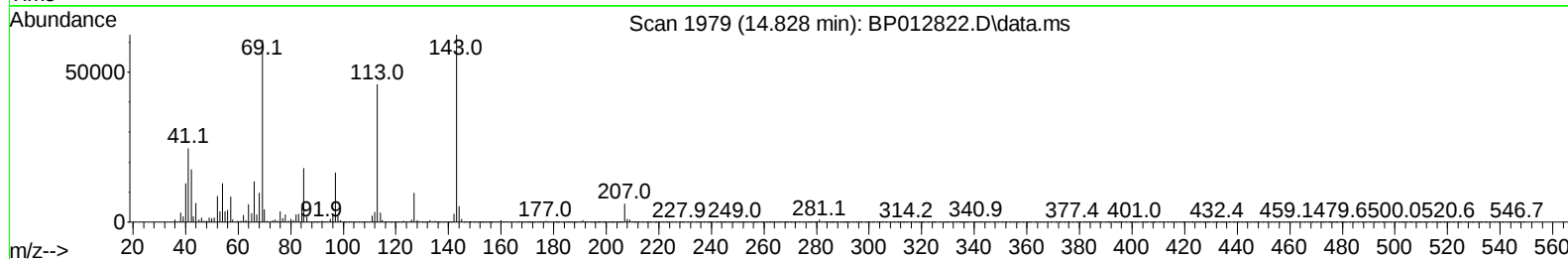
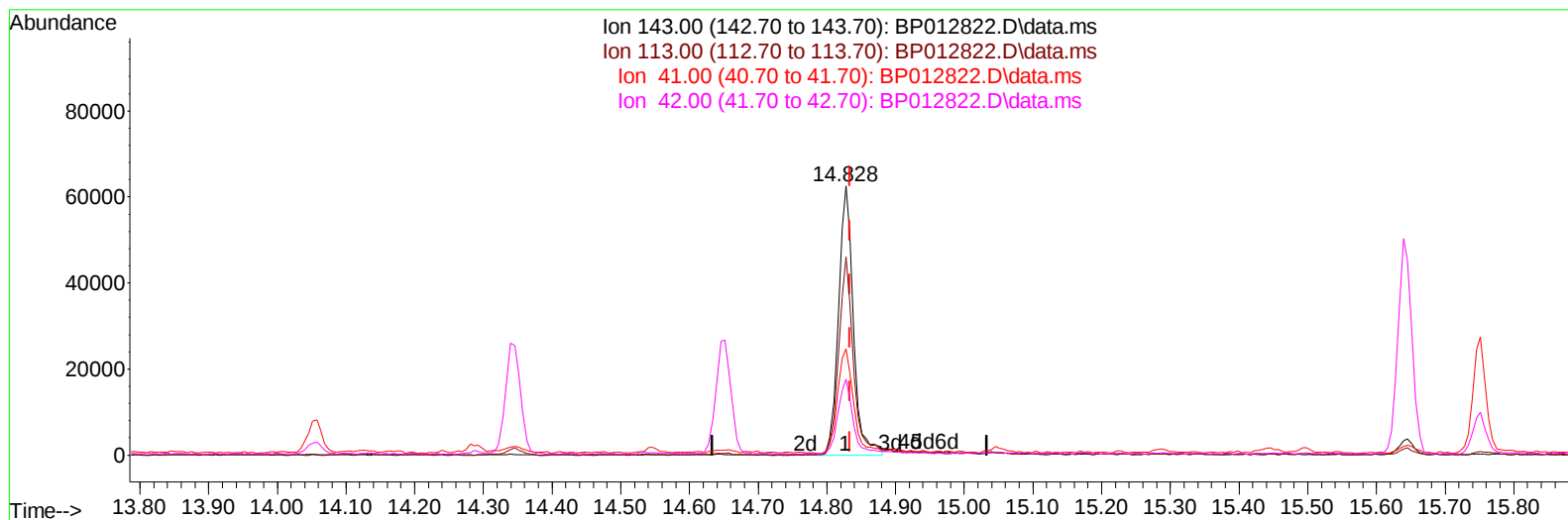
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TIC: BP012822.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.828min (-0.006) 4.94 ng/ul m

response 94067

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	65.70	73.64
41.00	37.00	39.44
42.00	24.70	28.22

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	553763	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	2451866	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1551724	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	2877072	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	1602765	20.000	ng/ul	0.00
88) Perylene-d12	24.004	264	1641993	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	64275	4.974	ng/uL	0.00
4) Pyridine-d5	3.828	84	136974	3.542	ng/ul	0.00
7) Phenol-d5	7.164	99	338992	7.450	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	917088	32.707	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	928483	26.778	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	571103	15.577	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	550416	36.678	ng/ul	0.00
24) 2-Nitrophenol-d4	9.905	143	585953	36.773	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	1081568	29.090	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	46364	0.911	ng/ul	0.00
46) Dimethylphthalate-d6	14.057	166	3813125	35.701	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	3878321	31.120	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	94067m	4.940	ng/ul	0.00
60) Fluorene-d10	15.645	176	3373022	36.060	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	436841	31.768	ng/ul	0.00
73) Anthracene-d10	17.504	188	4866776	38.428	ng/ul	0.00
81) Pyrene-d10	19.733	212	4632535	49.743	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.839	264	3227486	40.017	ng/ul	0.00

Target Compounds Qvalue

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

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