

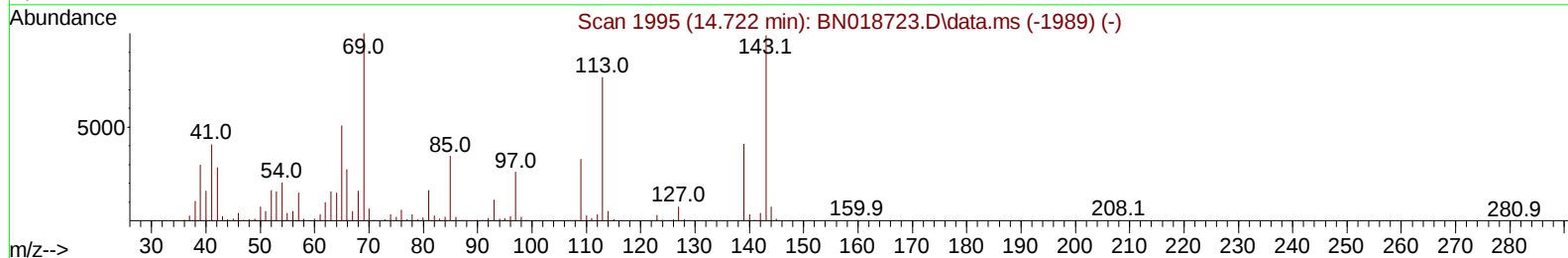
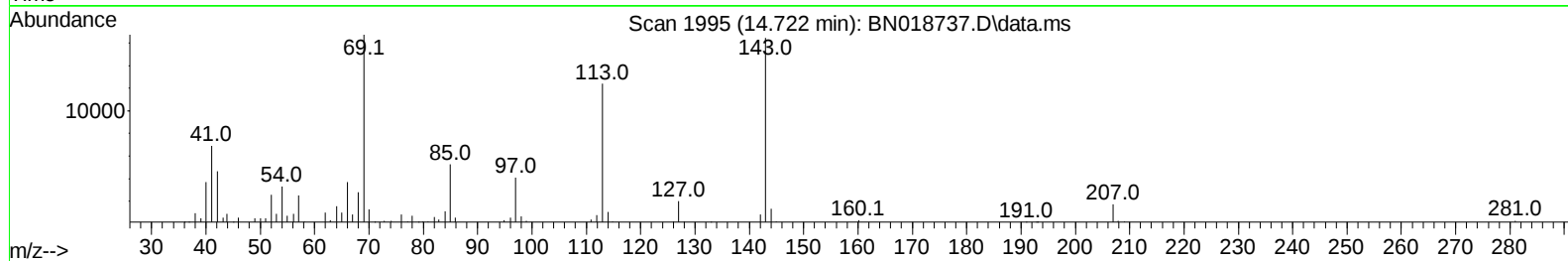
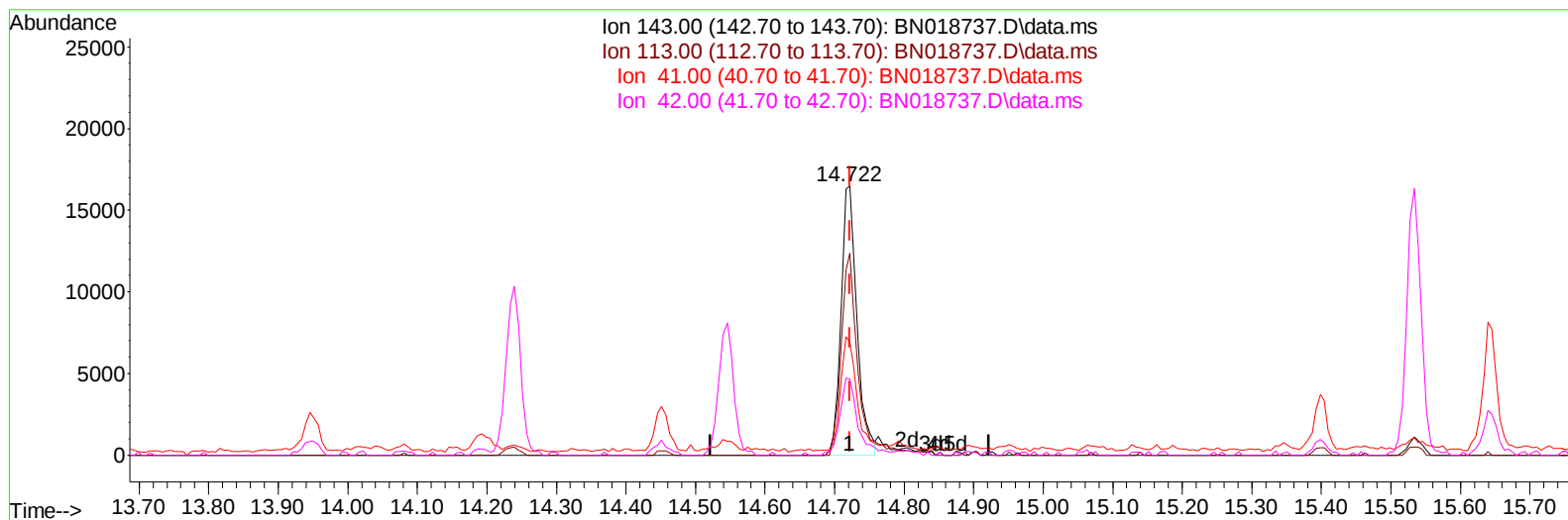
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
Data File : BN018737.D
Acq On : 25 Feb 2022 18:54
Operator : CG/JU
Sample : N1656-15
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGQ22

Manual IntegrationsAPPROVED

Quant Time: Feb 25 22:38:15 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 16 16:48:14 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/28/2022
Supervised By :mohammad ahmed 02/28/2022



TIC: BN018737.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.000) 5.71 ng/u1

response 26292

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	75.20
41.00	43.80	41.96
42.00	29.00	28.22

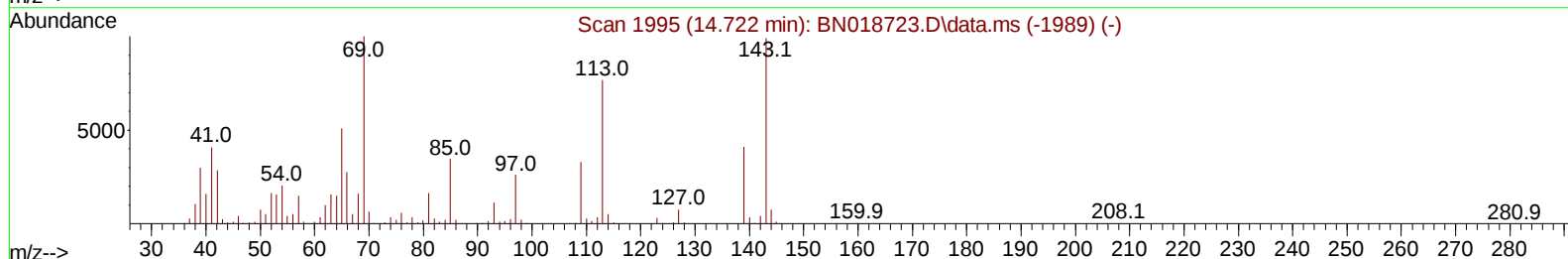
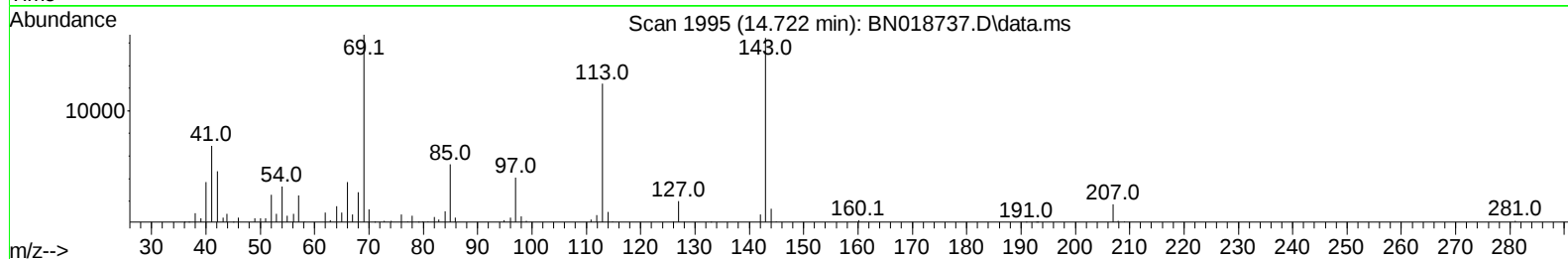
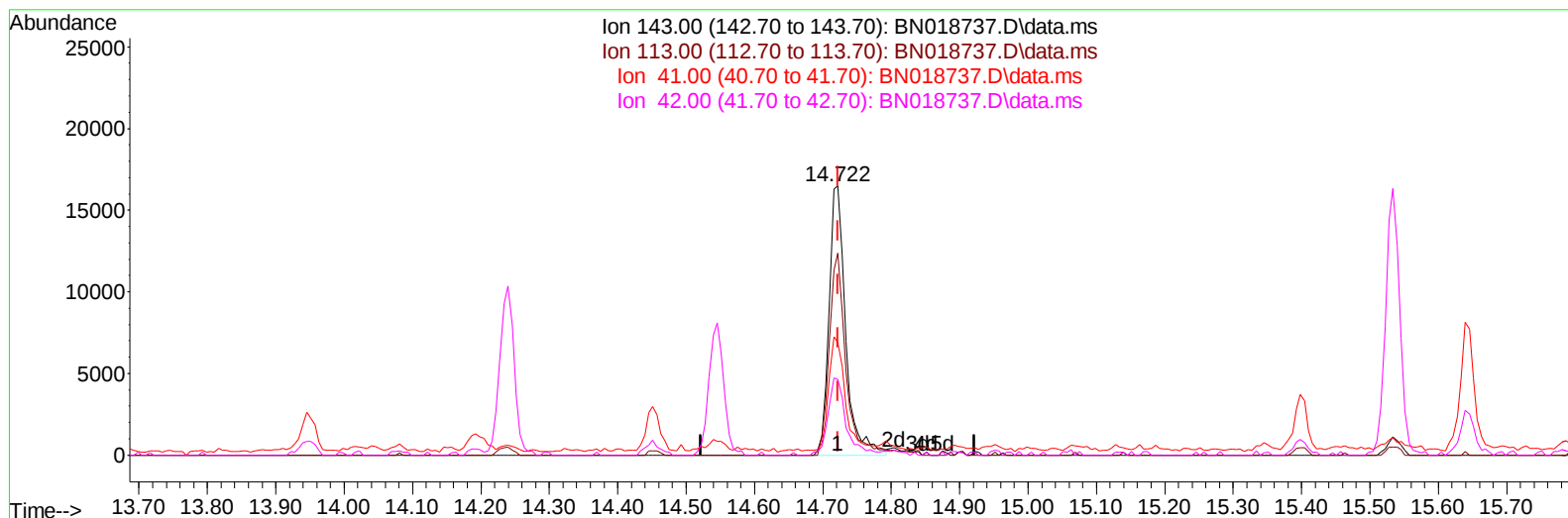
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TIC: BN018737.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.000) 5.97 ng/ul m

response 27512

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	75.20
41.00	43.80	41.96
42.00	29.00	28.22

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	177879	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	651536	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164	366045	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	709938	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	708363	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	744792	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	27414	5.809	ng/uL	0.00
4) Pyridine-d5	3.705	84	130615	10.237	ng/ul	0.00
7) Phenol-d5	7.063	99	122331	8.339	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	324650	35.655	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	337130	29.177	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	200604	17.906	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	182329	36.418	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	186962	34.869	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	319730	31.242	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	430953	31.516	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	995190	37.556	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1287264	37.496	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	27512m	5.975	ng/ul	0.00
60) Fluorene-d10	15.534	176	925351	40.448	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	120463	28.759	ng/ul	0.00
73) Anthracene-d10	17.381	188	1511615	44.498	ng/ul	0.00
81) Pyrene-d10	19.675	212	1750236	42.214	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1758793	46.304	ng/ul	0.00

Target Compounds Qvalue

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

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