

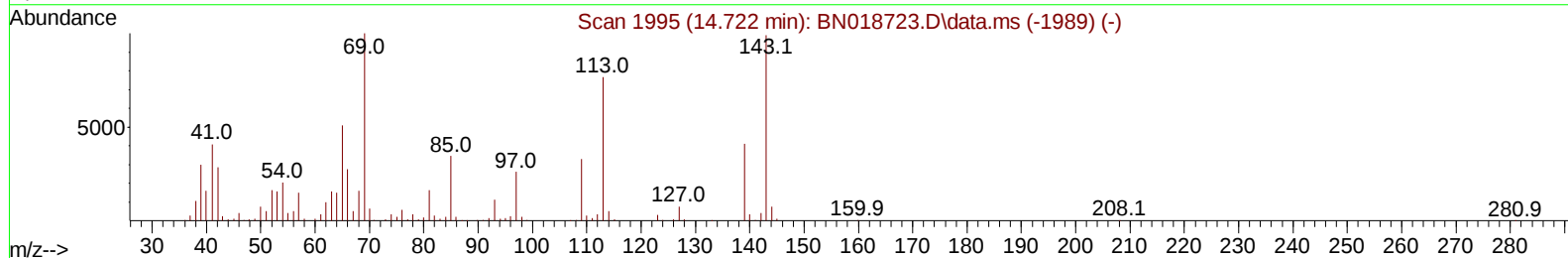
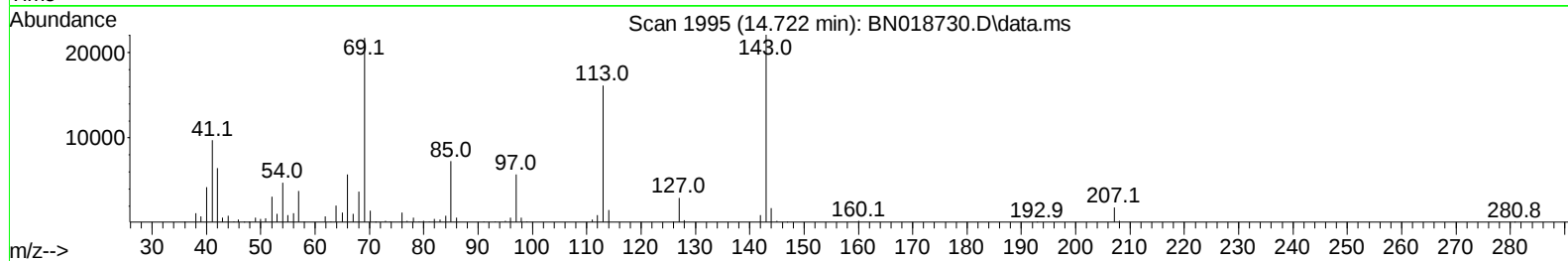
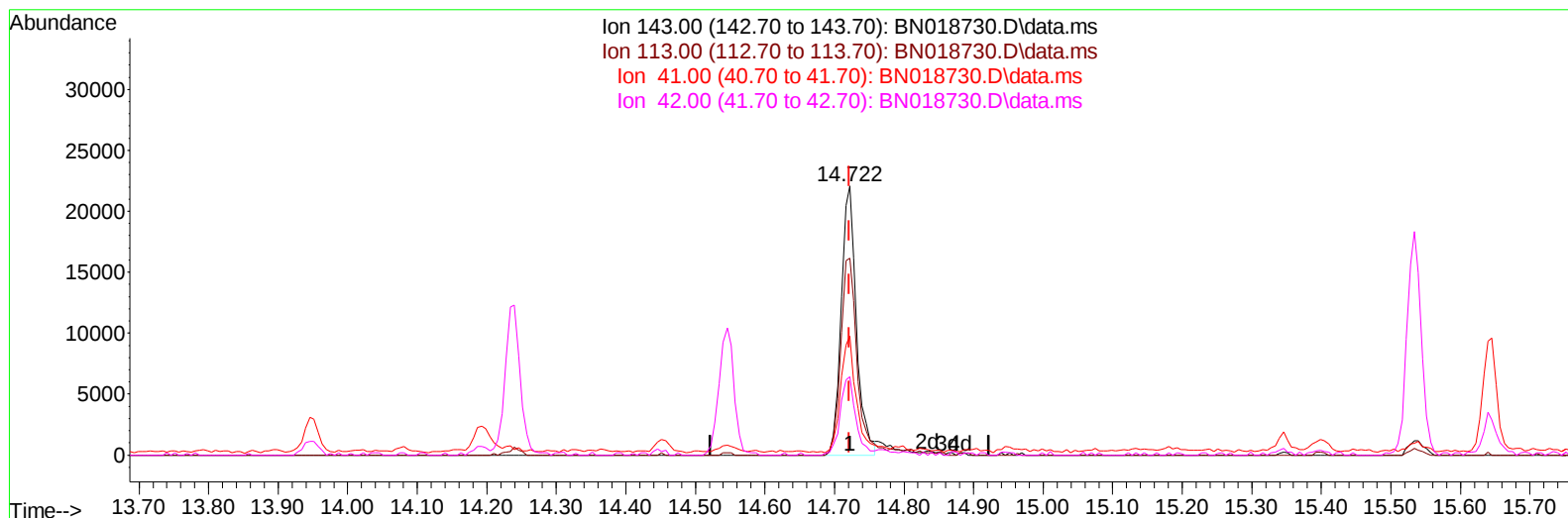
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
Data File : BN018730.D
Acq On : 25 Feb 2022 14:42
Operator : CG/JU
Sample : N1656-05
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGQ12

Manual IntegrationsAPPROVED

Quant Time: Feb 25 15:41:24 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: BN018730.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (+ 0.000) 5.38 ng/u1

response 33759

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	73.12
41.00	43.80	44.12
42.00	29.00	29.25

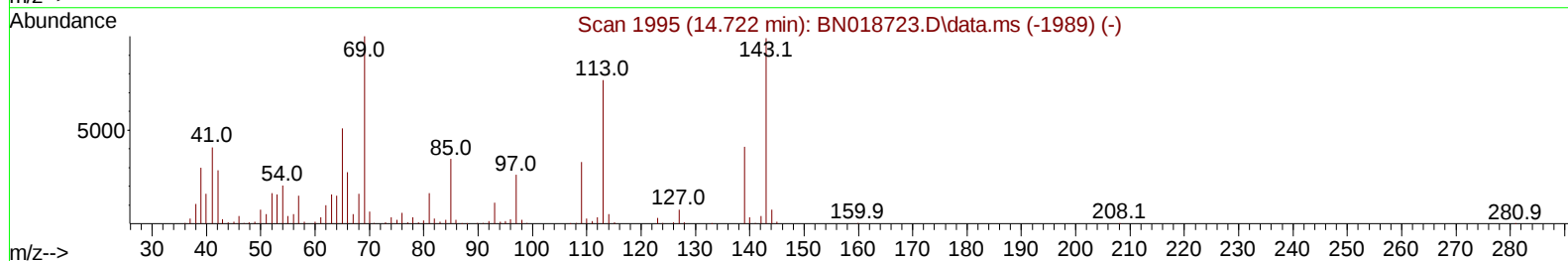
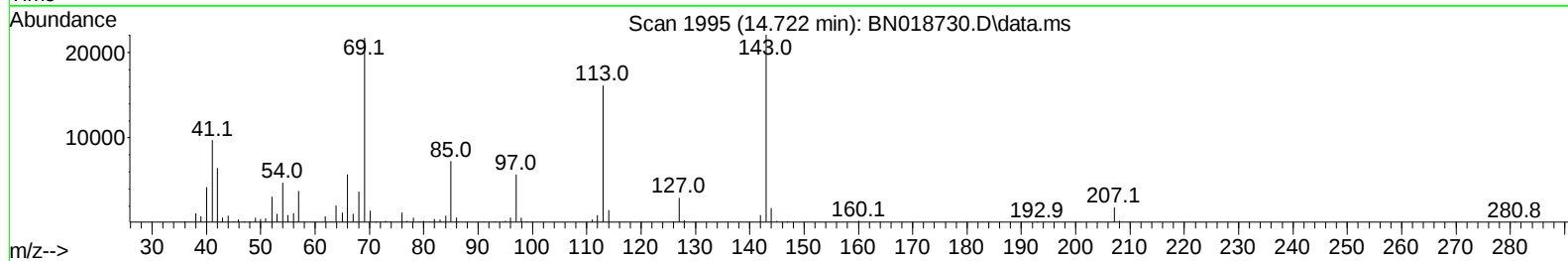
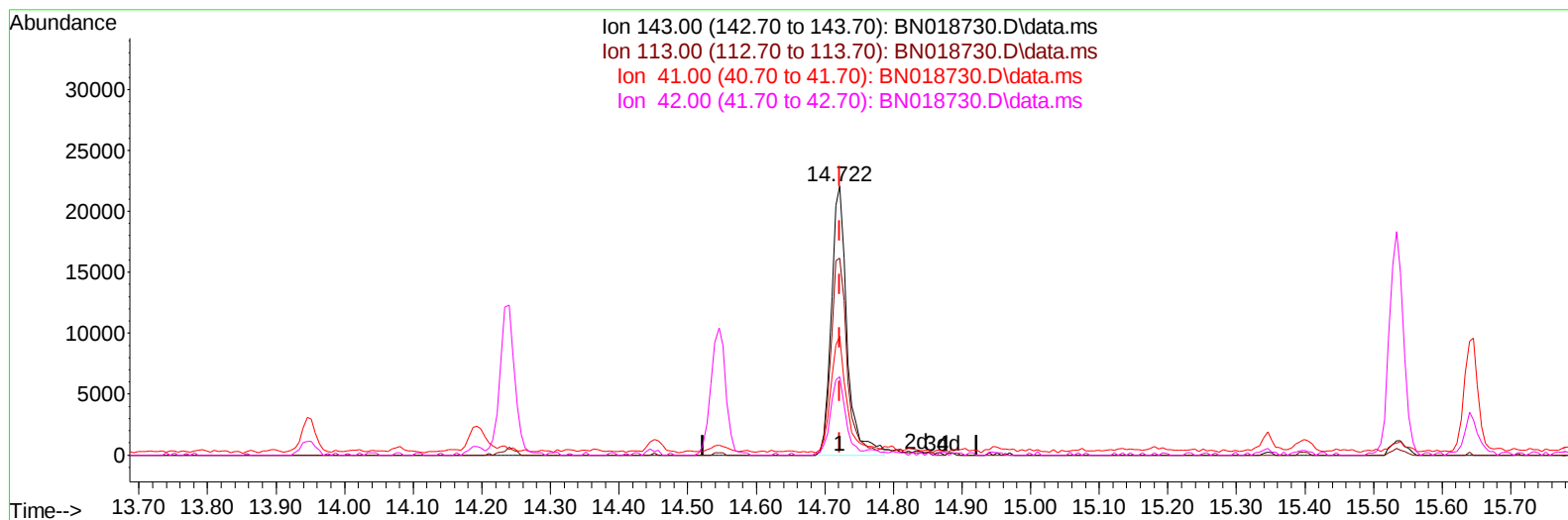
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TIC: BN018730.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (+ 0.000) 5.61 ng/ul m

response 35167

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	73.12
41.00	43.80	44.12
42.00	29.00	29.25

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	224629	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	880776	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.546	164	498596	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	946210	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	814534	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	796178	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	27041	4.537	ng/uL	0.00
4) Pyridine-d5	3.699	84	144371	8.960	ng/ul	0.00
7) Phenol-d5	7.064	99	144197	7.783	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	368769	32.072	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	374266	25.650	ng/ul	0.00
15) 4-Methylphenol-d8	8.611	113	243885	17.239	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	212043	31.330	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	220516	30.423	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	385156	27.839	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	482660	26.110	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	1209554	33.511	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160	1527790	32.672	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	35167m	5.607	ng/ul	0.00
60) Fluorene-d10	15.534	176	1023368	32.840	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	152995	27.405	ng/ul	0.00
73) Anthracene-d10	17.387	188	1637120	36.159	ng/ul	0.00
81) Pyrene-d10	19.675	212	1840242	38.599	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	1577877	38.860	ng/ul	0.00

Target Compounds Qvalue

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

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