

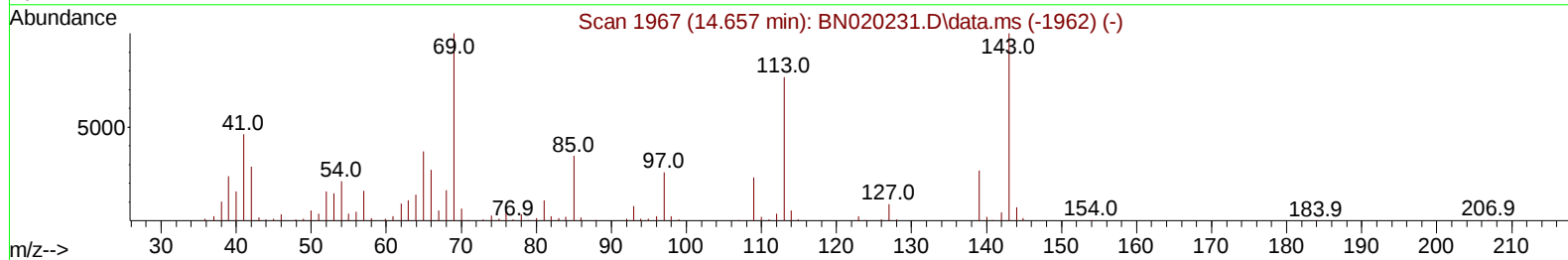
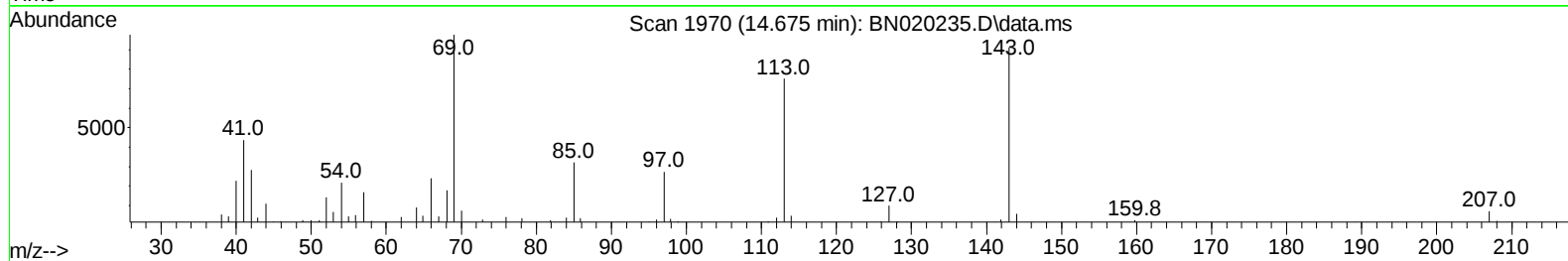
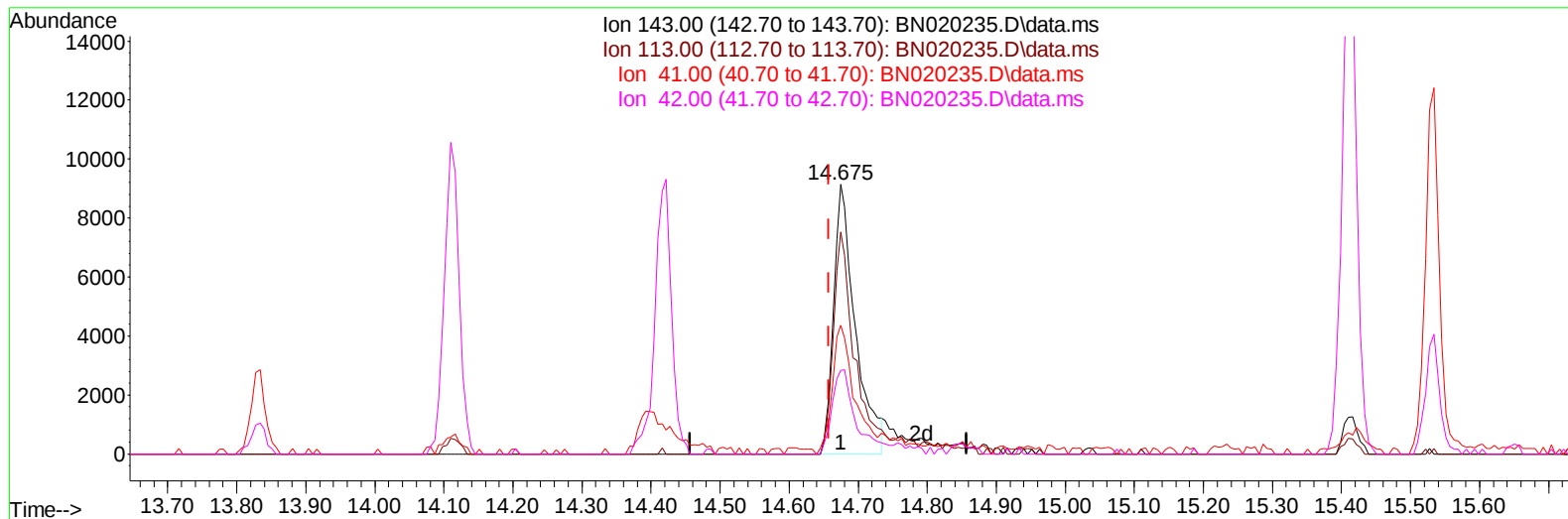
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061722\
Data File : BN020235.D
Acq On : 17 Jun 2022 12:30
Operator : CG/JU
Sample : N3313-09
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
YB7N1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/18/2022
Supervised By :mohammad ahmed 06/22/2022

Quant Time: Jun 17 22:56:19 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 22:53:31 2022
Response via : Initial Calibration



TIC: BN020235.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.017) 3.02 ng/u1

response 19956

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	82.31
41.00	39.70	47.78#
42.00	27.00	30.91

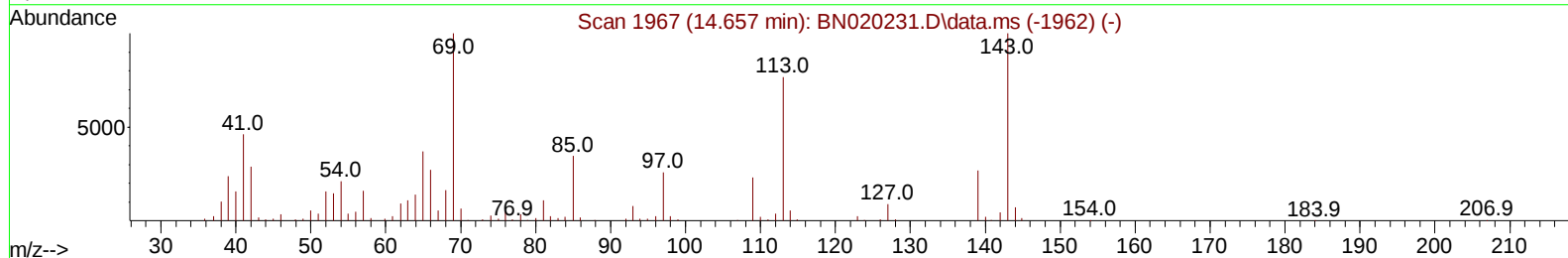
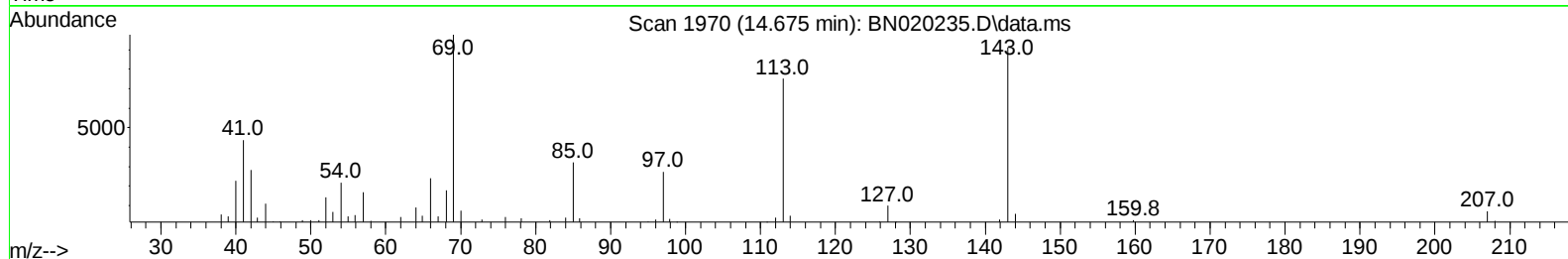
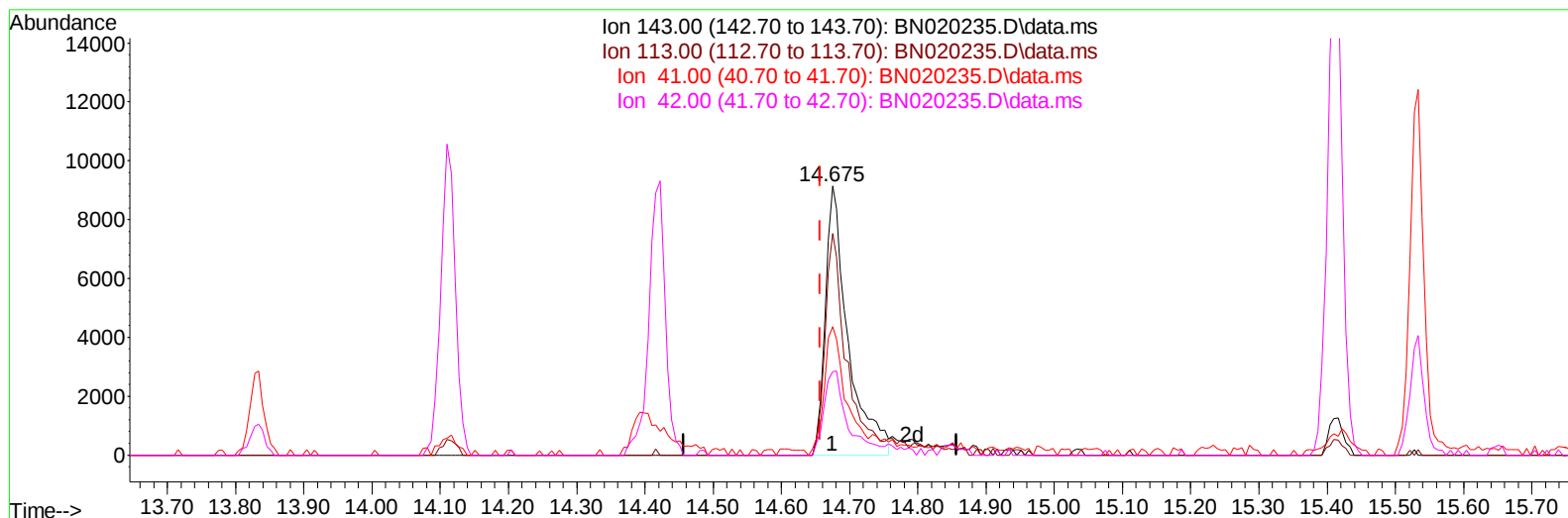
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TIC: BN020235.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.017) 3.20 ng/ul m

response 21145

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	82.31
41.00	39.70	47.78#
42.00	27.00	30.91

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	129139	20.000	ng/ul	0.00
20) Naphthalene-d8	10.581	136	650718	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	453772	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.163	188	963410	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	855542	20.000	ng/ul	0.00
88) Perylene-d12	23.662	264	700859	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	15177	5.090	ng/uL	0.00
4) Pyridine-d5	3.664	84	82630	10.139	ng/ul	0.00
7) Phenol-d5	6.987	99	67896	6.354	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	217270	32.543	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	222029	25.572	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	150960	16.329	ng/ul	0.00
21) Nitrobenzene-d5	8.946	128	150889	30.479	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	170146	32.385	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	279174	30.438	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	404492	27.233	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1085371	32.873	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1254136	32.754	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	21145m	3.196	ng/ul	0.02
60) Fluorene-d10	15.410	176	968153	35.435	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.534	200	168918	30.268	ng/ul	0.00
73) Anthracene-d10	17.263	188	1611950	38.338	ng/ul	0.00
81) Pyrene-d10	19.557	212	1784133	38.467	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	1324563	38.627	ng/ul	0.00

Target Compounds Qvalue

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

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