

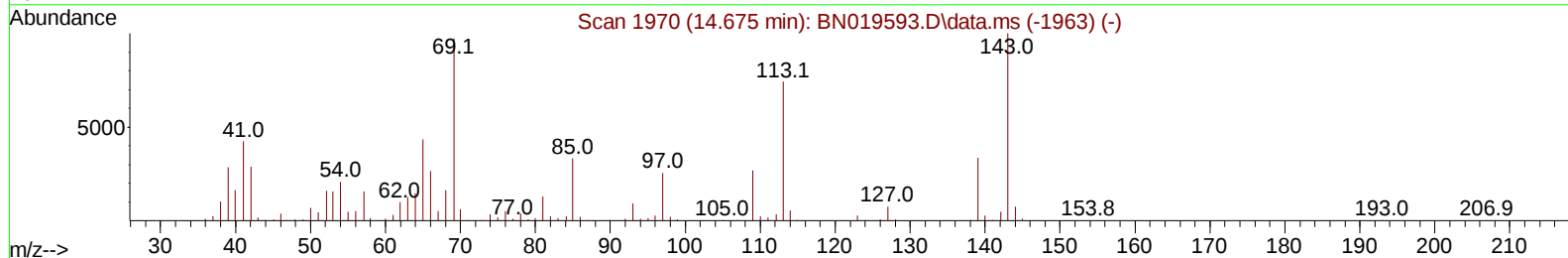
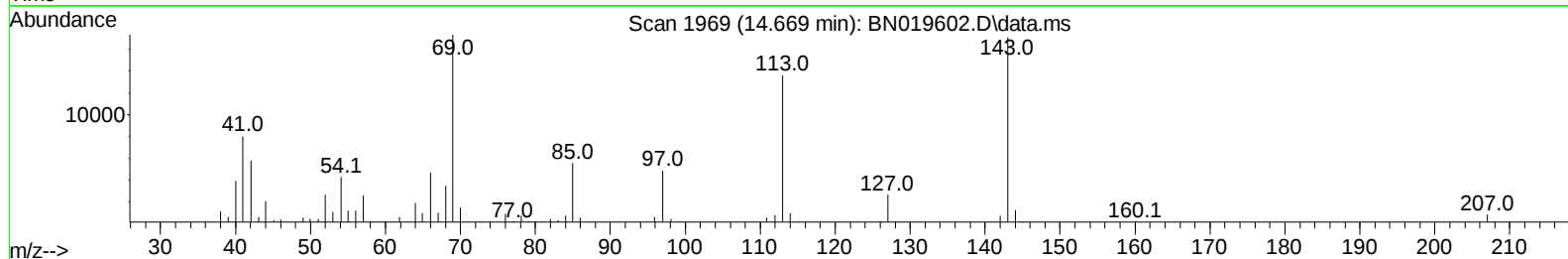
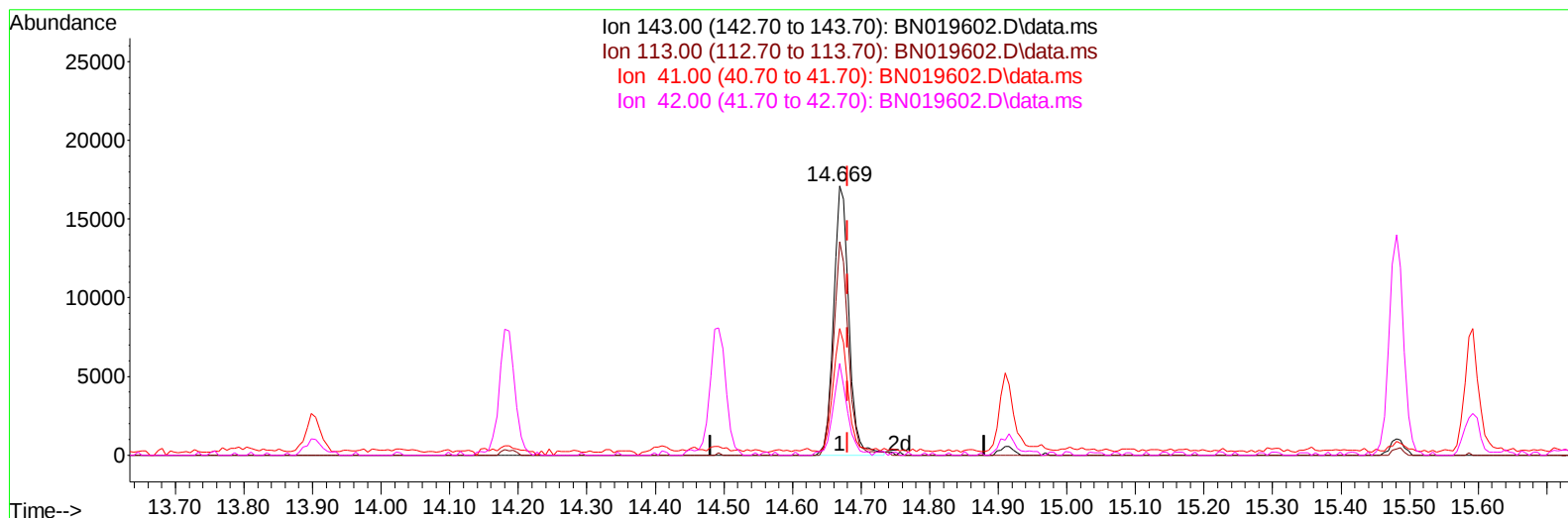
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043022\
Data File : BN019602.D
Acq On : 29 Apr 2022 22:21
Operator : CG/JU
Sample : N2612-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH0B2

Manual IntegrationsAPPROVED

Quant Time: Apr 30 03:07:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 27 13:41:44 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/02/2022
Supervised By :Yogesh Patel 05/05/2022



TIC: BN019602.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.012) 5.06 ng/u1

response 26335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	79.45
41.00	39.70	47.06
42.00	27.00	34.04#

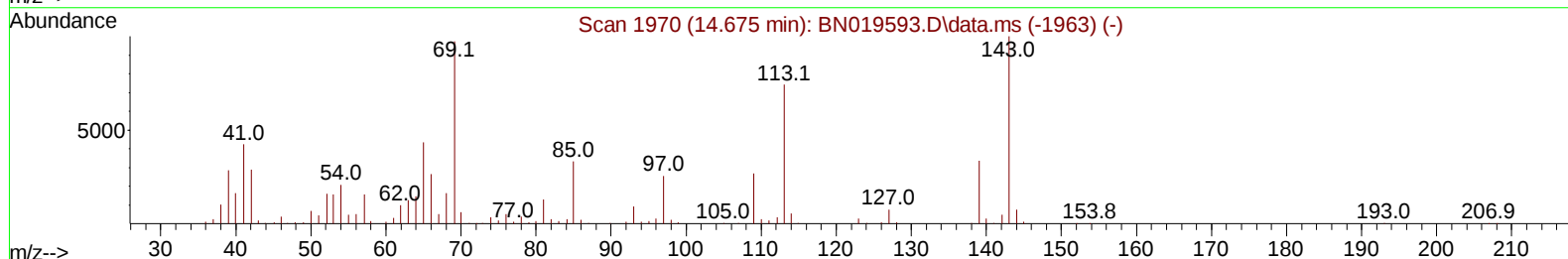
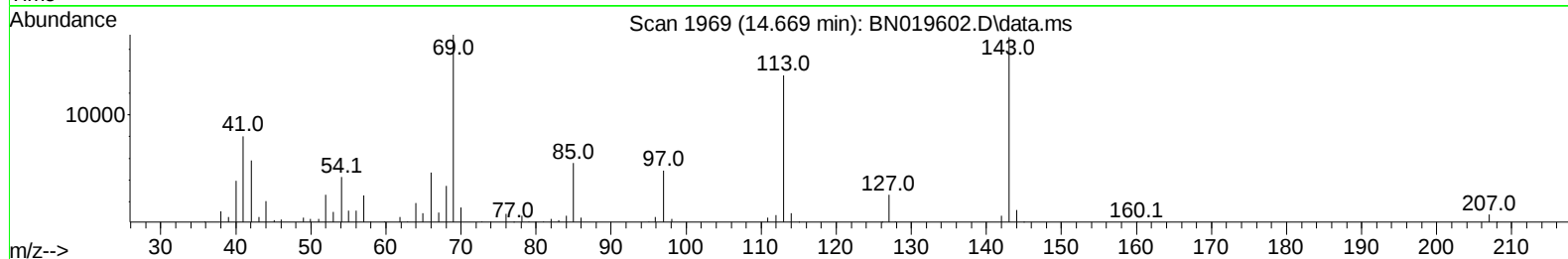
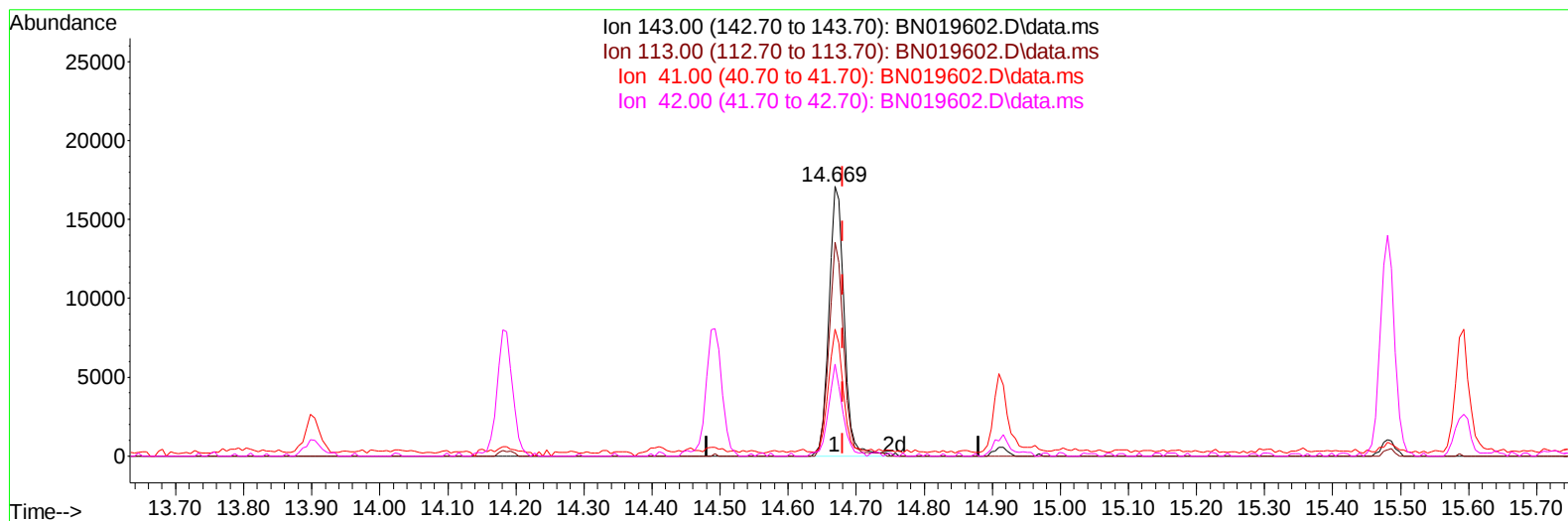
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TIC: BN019602.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.012) 5.09 ng/u1 m

response 26473

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	79.45
41.00	39.70	47.06
42.00	27.00	34.04#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	151747	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	642338	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	389395	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	811358	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	754017	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	768077	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	15443	3.840	ng/uL	0.00
4) Pyridine-d5	3.752	84	77424	7.604	ng/ul	0.00
7) Phenol-d5	7.034	99	70172	5.336	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	213738	28.021	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.404	132	222575	21.338	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	142066	14.102	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	144789	30.916	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	140529	30.448	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	236977	25.650	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	367982	26.238	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	868229	32.463	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	1027309	28.408	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	26473m	5.089	ng/ul	-0.01
60) Fluorene-d10	15.481	176	763028	32.318	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	102024	24.631	ng/ul	0.00
73) Anthracene-d10	17.328	188	1214752	31.792	ng/ul	0.00
81) Pyrene-d10	19.622	212	1444336	33.317	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1327111	33.800	ng/ul	0.00

Target Compounds Qvalue

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

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