

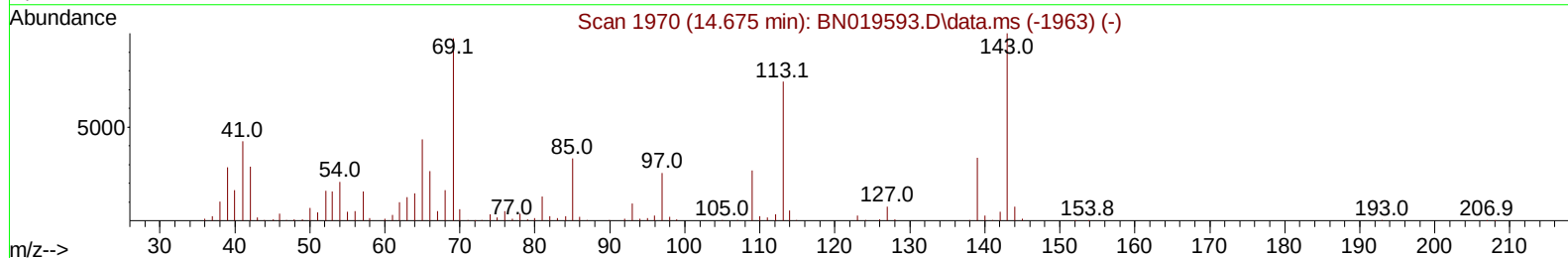
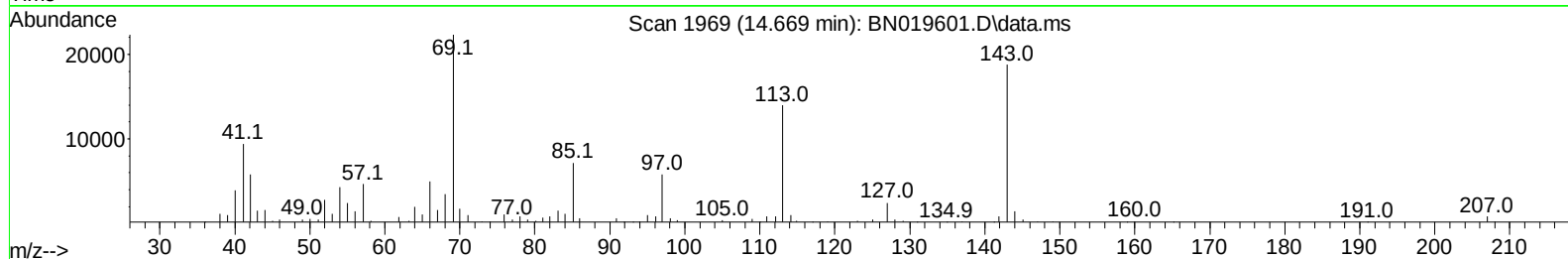
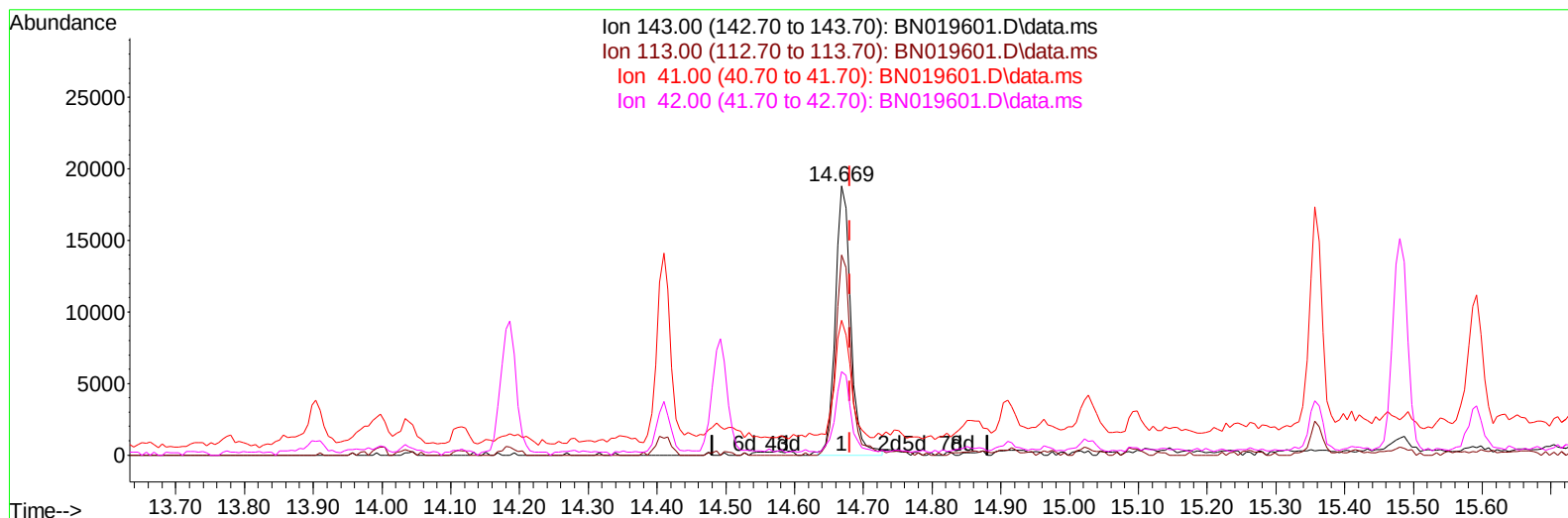
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN043022\
Data File : BN019601.D
Acq On : 29 Apr 2022 21:45
Operator : CG/JU
Sample : N2612-04
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH099

Manual IntegrationsAPPROVED

Quant Time: Apr 30 03:07:18 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: BN019601.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.012) 5.95 ng/u1

response 29578

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	74.36
41.00	39.70	50.10#
42.00	27.00	31.02

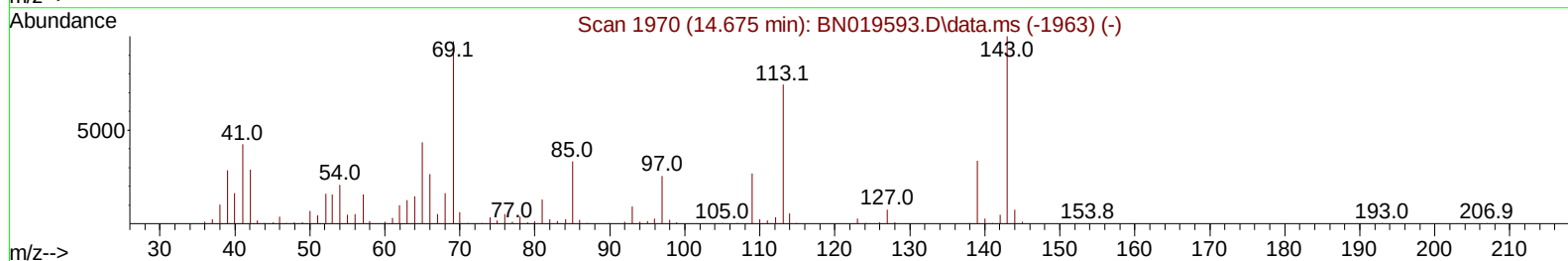
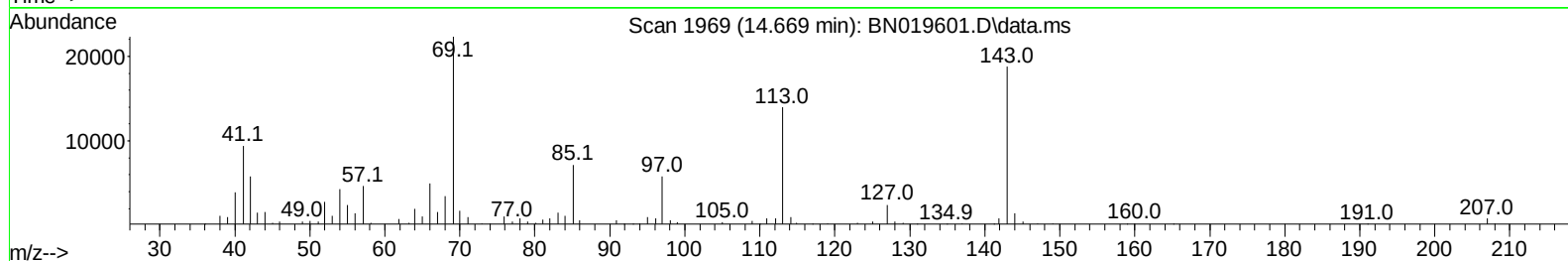
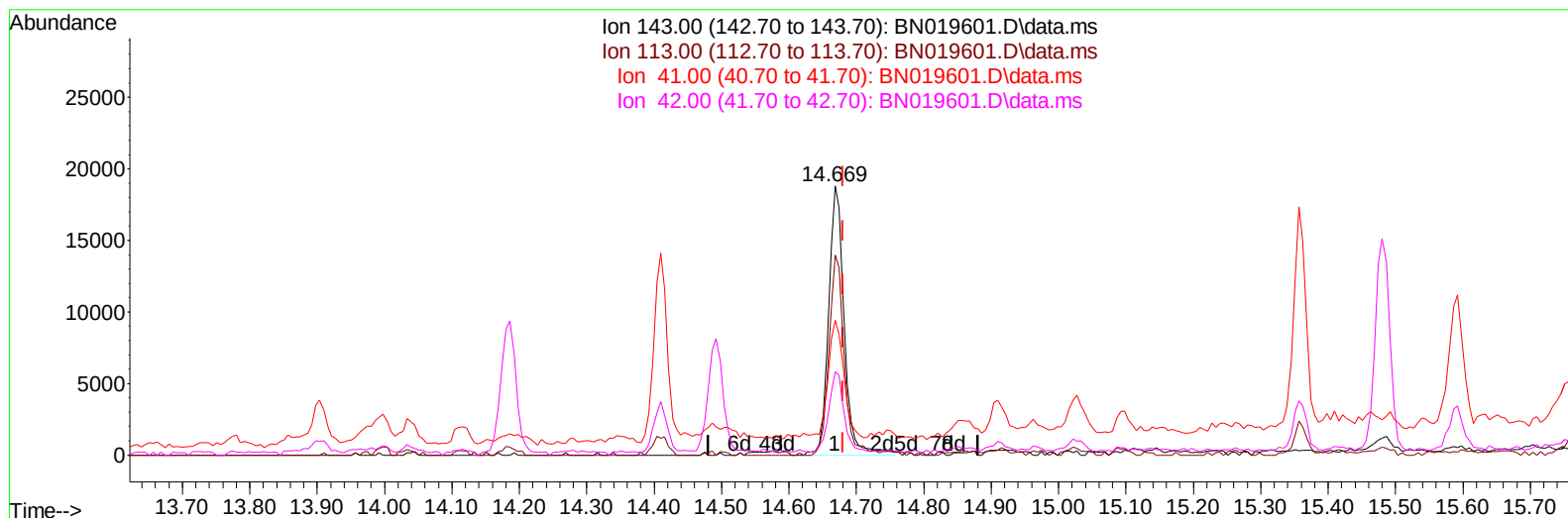
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TIC: BN019601.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.012) 6.08 ng/ul m

response 30248

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	74.36
41.00	39.70	50.10#
42.00	27.00	31.02

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	156819	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	634856	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	372354	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	747942	20.000	ng/ul	-0.01
79) Chrysene-d12	21.415	240	701386	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	714271	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	13618	3.277	ng/uL	0.00
4) Pyridine-d5	3.751	84	61911	5.883	ng/ul	0.00
7) Phenol-d5	7.034	99	75932	5.587	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.198	67	245043	31.087	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.404	132	258896	24.018	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	153196	14.715	ng/ul	-0.01
21) Nitrobenzene-d5	9.022	128	161790	34.953	ng/ul	0.00
24) 2-Nitrophenol-d4	9.745	143	163196	35.776	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	260536	28.532	ng/ul	0.00
31) 4-Chloroaniline-d4	10.792	131	343412	24.774	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	948030	37.069	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	1123172	32.481	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	30248m	6.081	ng/ul	-0.01
60) Fluorene-d10	15.480	176	815943	36.141	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	117534	30.782	ng/ul	0.00
73) Anthracene-d10	17.327	188	1330251	37.766	ng/ul	0.00
81) Pyrene-d10	19.621	212	1536805	38.111	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1469389	40.243	ng/ul	0.00

Target Compounds Qvalue

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

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