

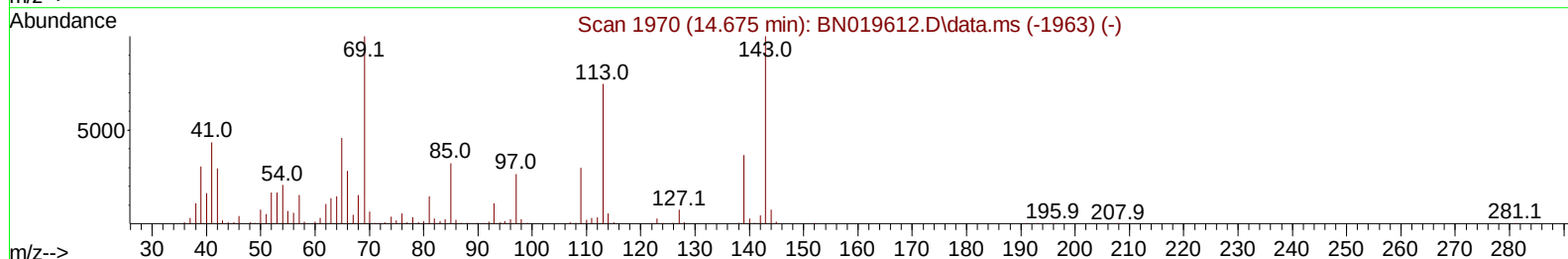
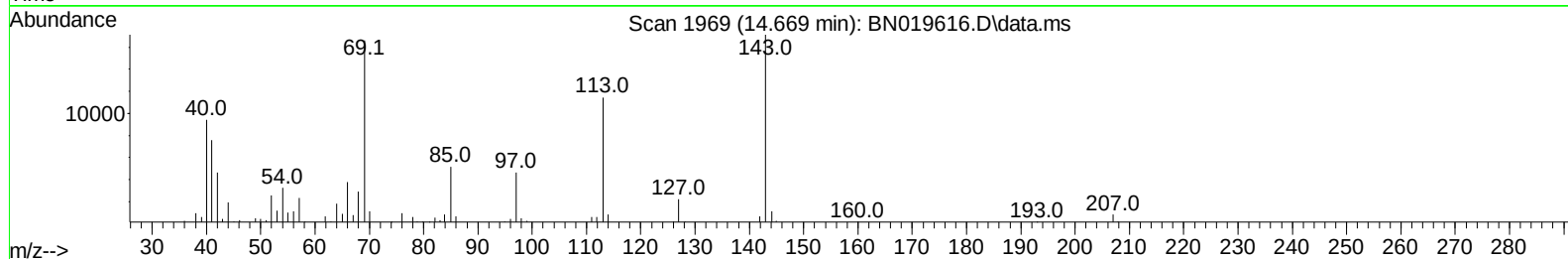
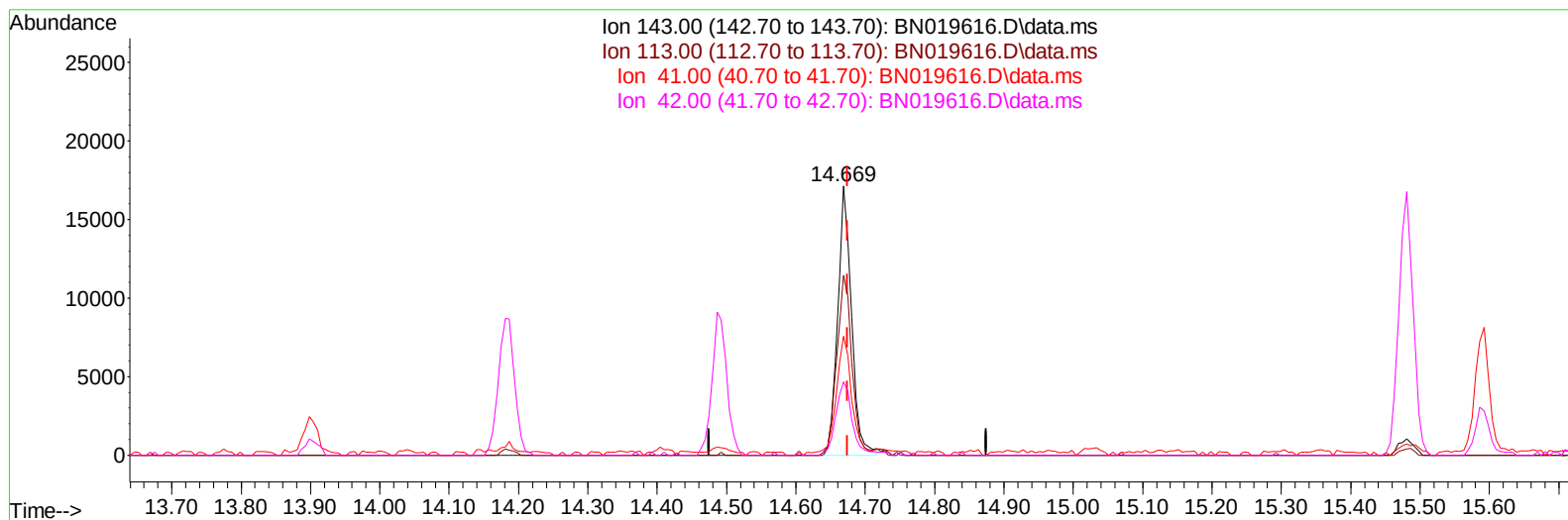
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019616.D
Acq On : 02 May 2022 13:12
Operator : CG/JU
Sample : N2612-22
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH0D3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :Yogesh Patel 05/09/2022

Quant Time: May 03 03:44:08 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 03 03:31:44 2022
Response via : Initial Calibration



TIC: BN019616.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 4.47 ng/ul

response 23450

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	66.80
41.00	39.70	44.11
42.00	27.00	27.22

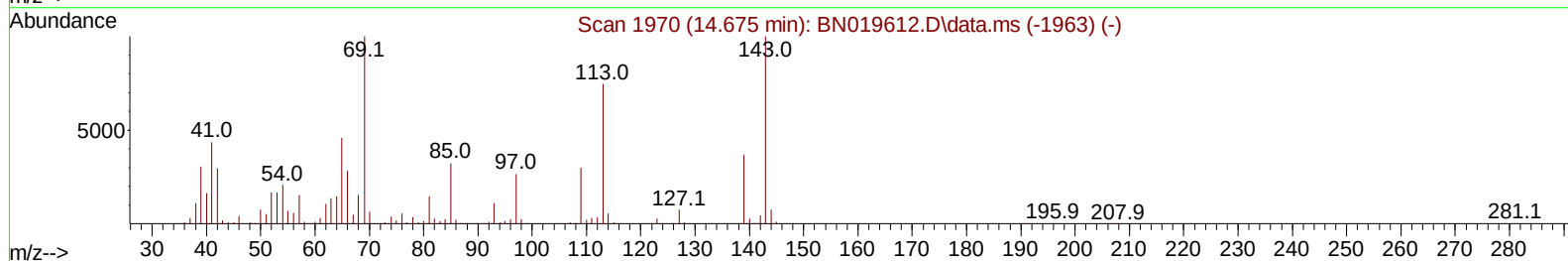
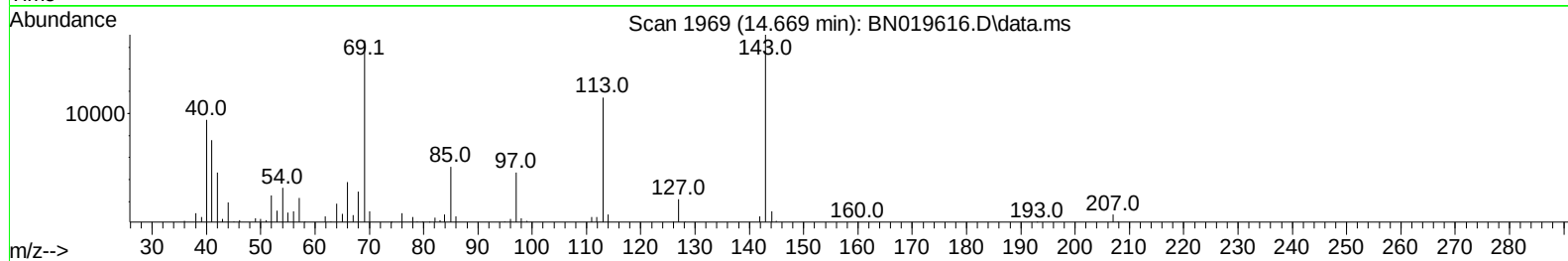
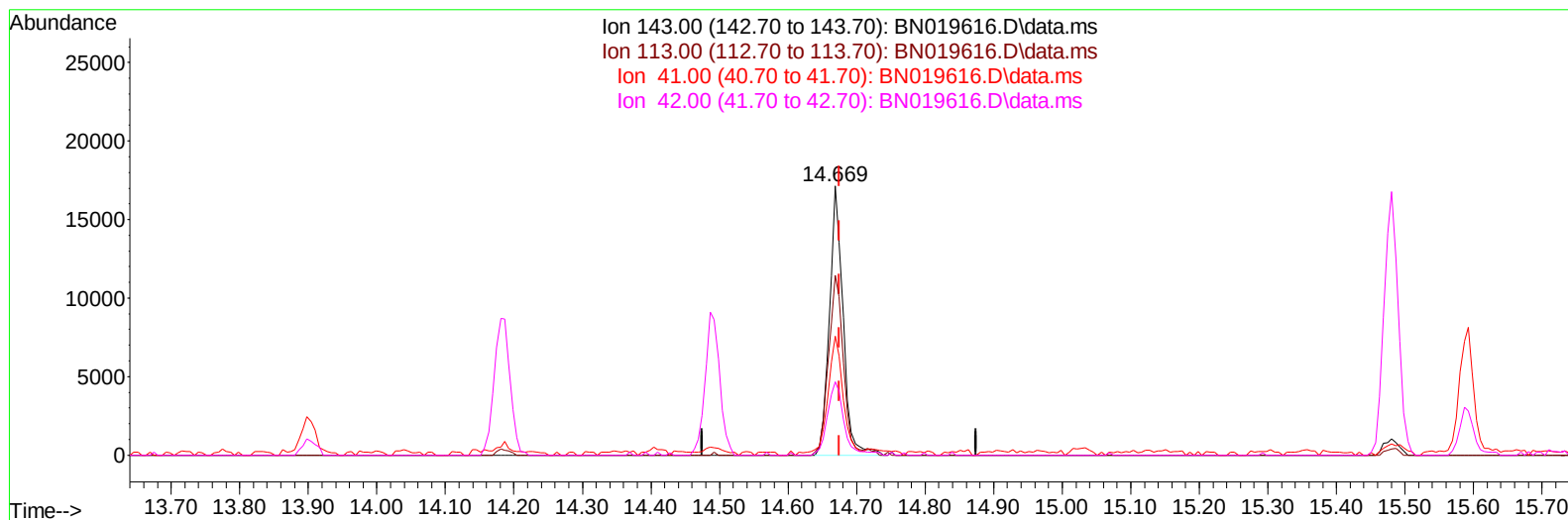
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TIC: BN019616.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 4.53 ng/ul m

response 23757

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	66.80
41.00	39.70	44.11
42.00	27.00	27.22

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	157269	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	666701	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	392962	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	789726	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	755303	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	767503	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	21216	5.091	ng/uL	0.00
4) Pyridine-d5	3.752	84	33023	3.129	ng/ul	0.00
7) Phenol-d5	7.034	99	69584	5.106	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	257503	32.574	ng/ul	0.00
11) 2-Chlorophenol-d4	7.404	132	248568	22.993	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	144197	13.811	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	160845	33.089	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	145876	30.451	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	259907	27.104	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	317419	21.805	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	919657	34.073	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	1124789	30.822	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	23757m	4.526	ng/ul	0.00
60) Fluorene-d10	15.481	176	836024	35.088	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	109989	27.281	ng/ul	0.00
73) Anthracene-d10	17.328	188	1481778	39.843	ng/ul	0.00
81) Pyrene-d10	19.616	212	1672984	38.526	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.615	264	1597875	40.726	ng/ul	0.00

Target Compounds Qvalue

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

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