

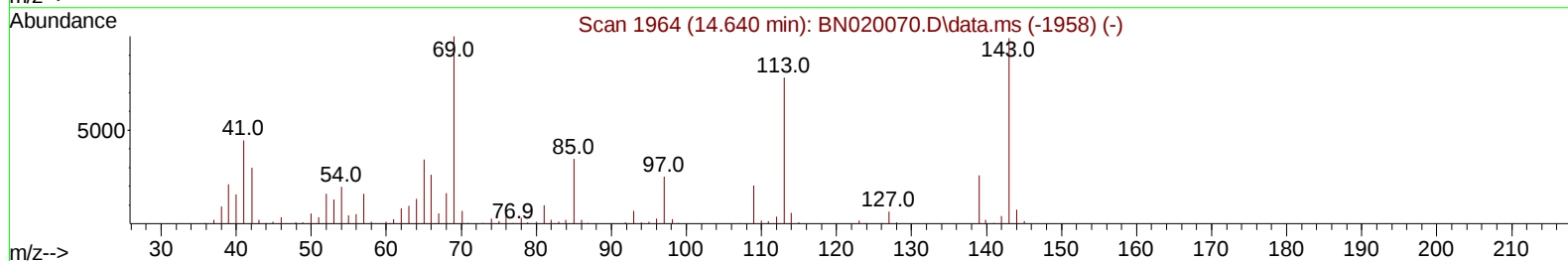
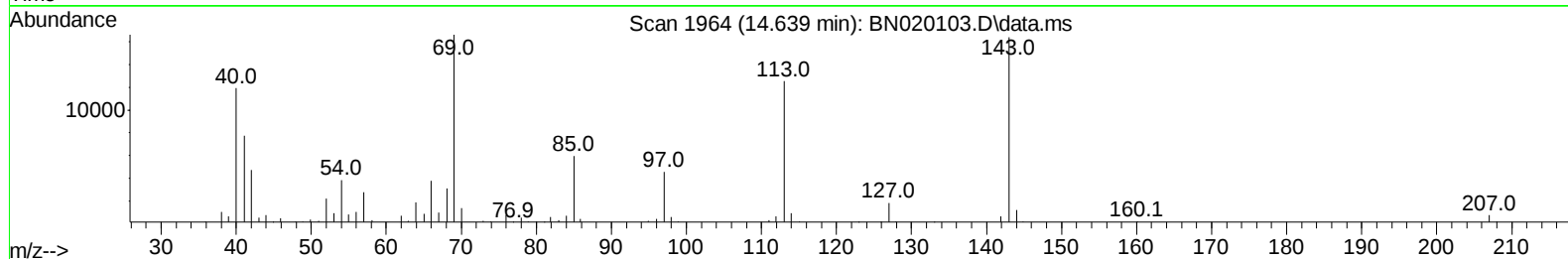
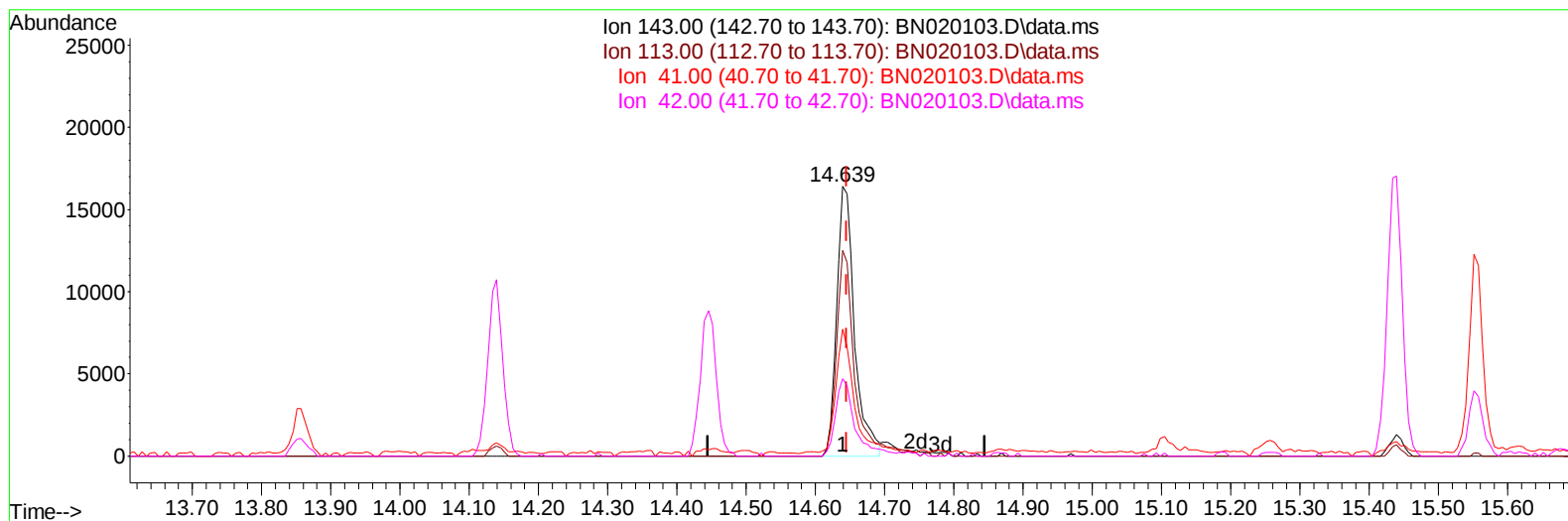
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061122\
Data File : BN020103.D
Acq On : 10 Jun 2022 16:43
Operator : CG/JU
Sample : N3250-15
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EXYT0

Manual IntegrationsAPPROVED

Quant Time: Jun 10 22:46:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 06 15:14:59 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/13/2022
Supervised By :mohammad ahmed 06/14/2022



TIC: BN020103.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.639min (-0.006) 4.66 ng/u1

response 29294

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	76.29
41.00	39.70	47.15
42.00	27.00	28.67

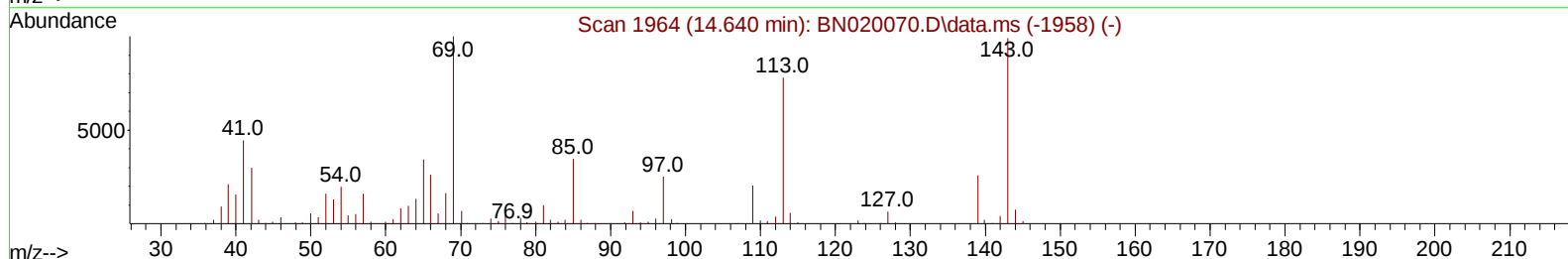
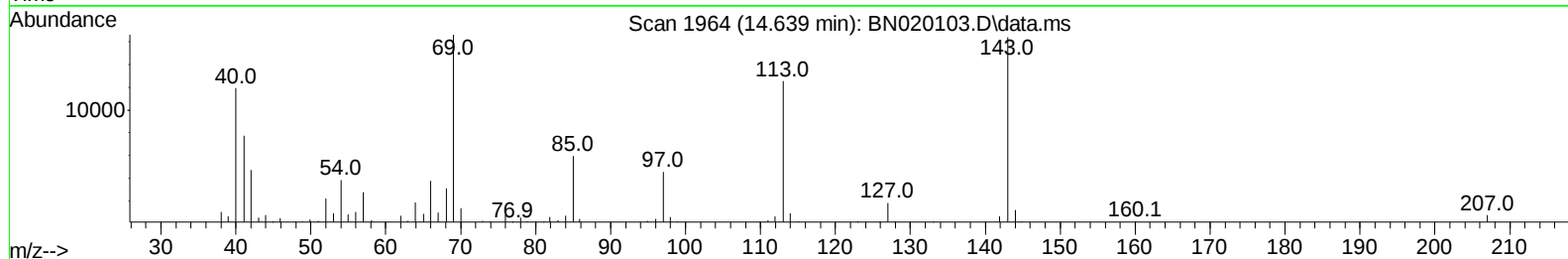
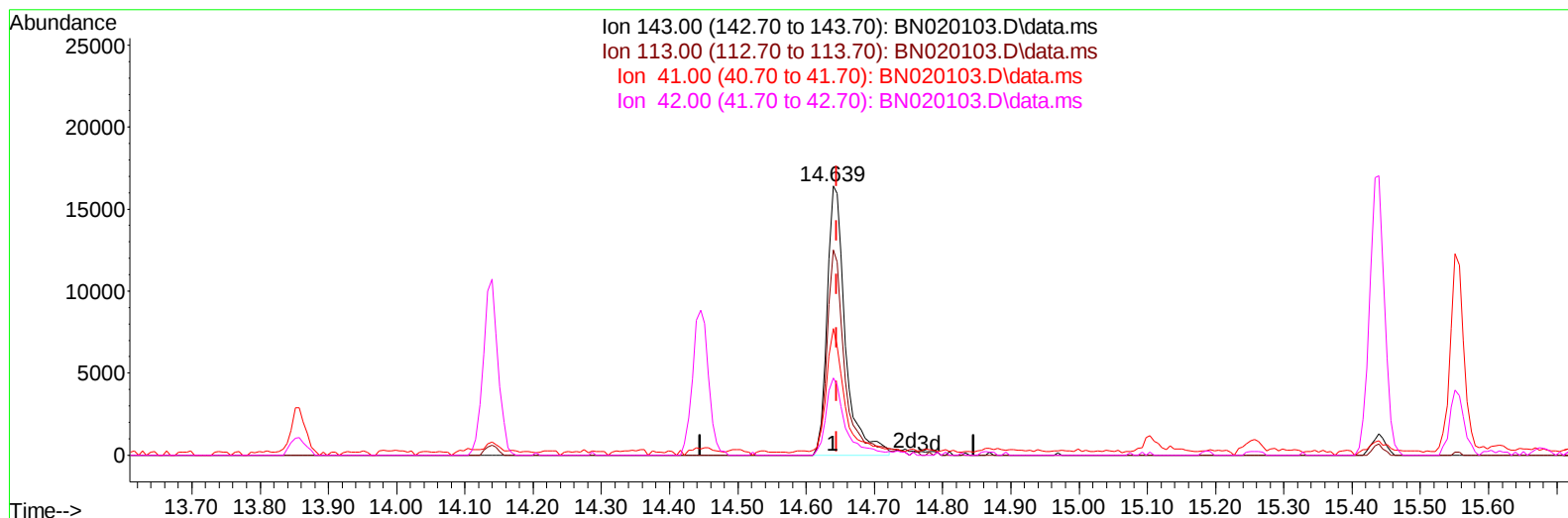
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TIC: BN020103.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.639min (-0.006) 4.84 ng/ul m

response 30419

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	76.29
41.00	39.70	47.15
42.00	27.00	28.67

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	137125	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	651312	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	431009	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	907464	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	825239	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	754516	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	13753	4.344	ng/uL	0.00
4) Pyridine-d5	3.693	84	73127	8.451	ng/ul	0.00
7) Phenol-d5	6.987	99	71019	6.259	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	244026	34.422	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	244556	26.526	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	155368	15.827	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	168539	34.014	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	178104	33.869	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	274720	29.925	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	370030	24.890	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	1095272	34.925	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	1259116	34.621	ng/ul	0.00
54) 4-Nitrophenol-d4	14.639	143	30419m	4.840	ng/ul	0.00
60) Fluorene-d10	15.439	176	927206	35.729	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	164113	31.220	ng/ul	0.00
73) Anthracene-d10	17.286	188	1538469	38.846	ng/ul	0.00
81) Pyrene-d10	19.580	212	1730718	38.686	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.557	264	1468924	39.790	ng/ul	0.00

Target Compounds Qvalue

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

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