

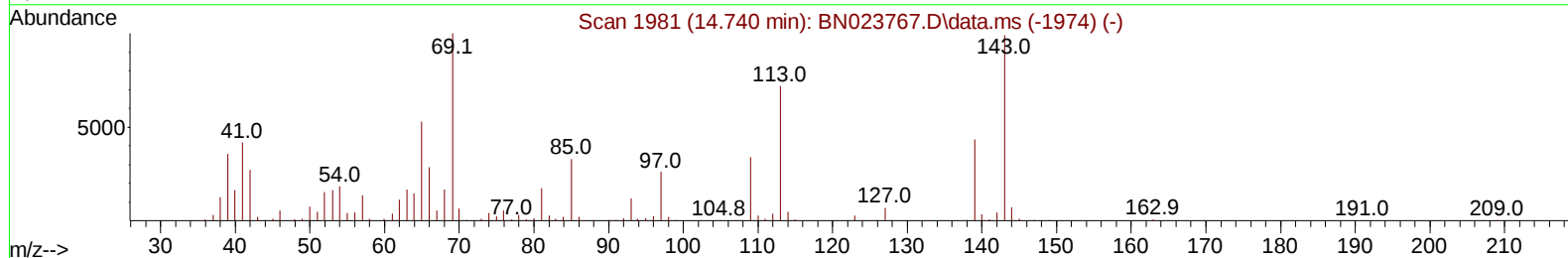
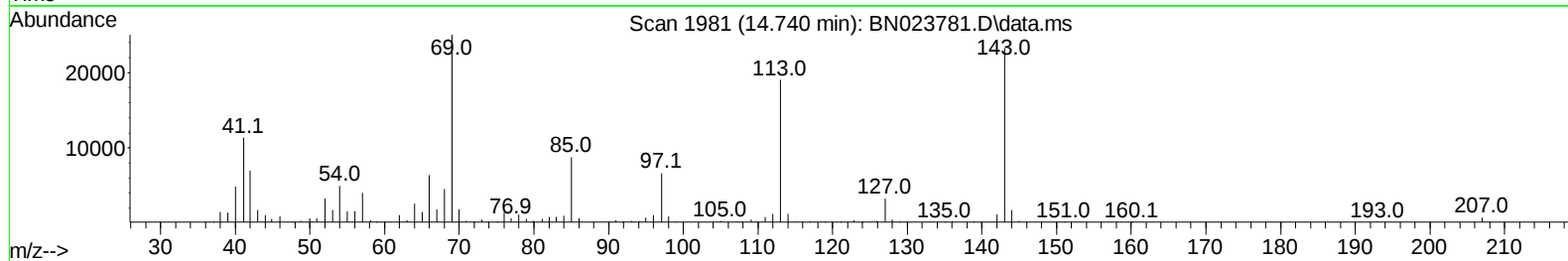
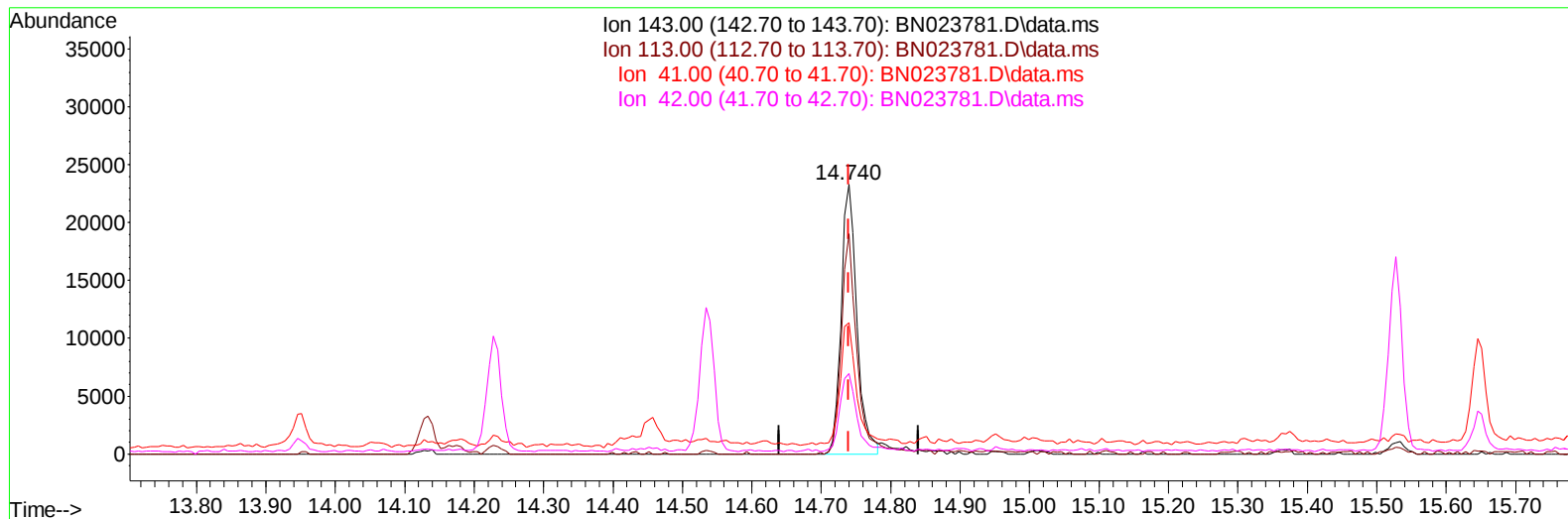
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN012323\
Data File : BN023781.D
Acq On : 24 Jan 2023 17:36
Operator : CG/JU
Sample : 01226-15
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EX9G2

Manual IntegrationsAPPROVED

Quant Time: Jan 24 23:58:27 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010523.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 24 23:52:16 2023
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/25/2023
Supervised By :Yogesh Patel 01/25/2023



TIC: BN023781.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (-0.000) 4.95 ng/u1

response 36100

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	81.63
41.00	40.90	48.81
42.00	27.80	29.89

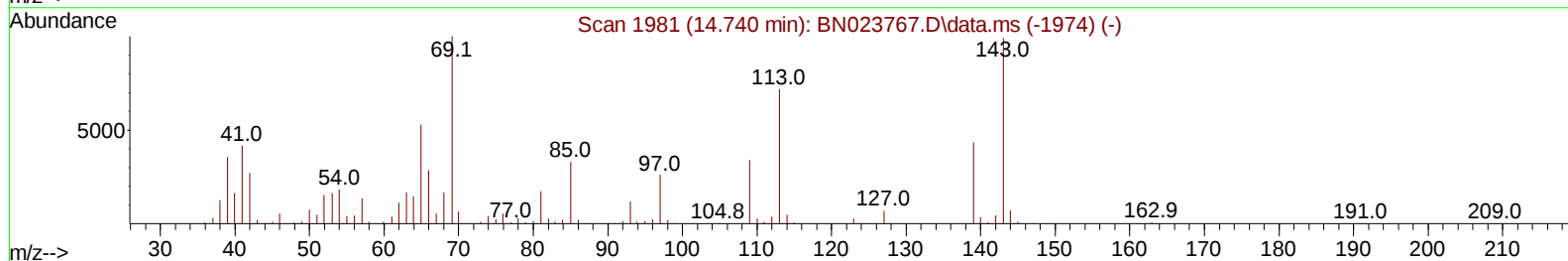
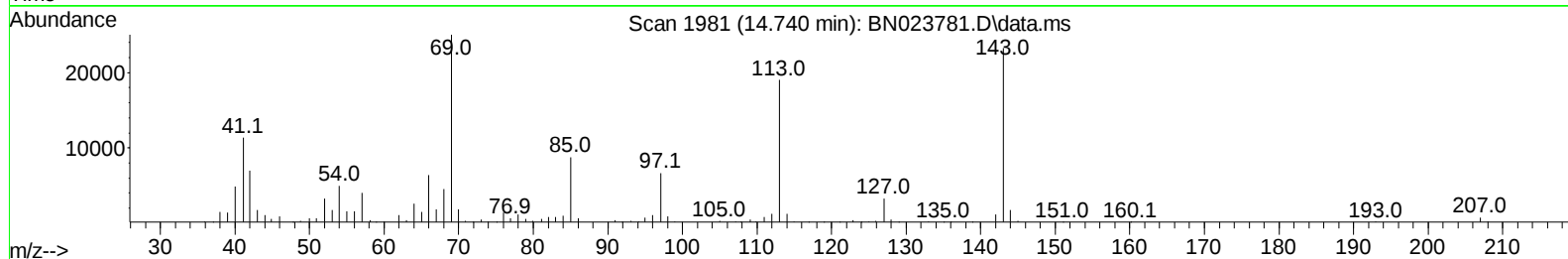
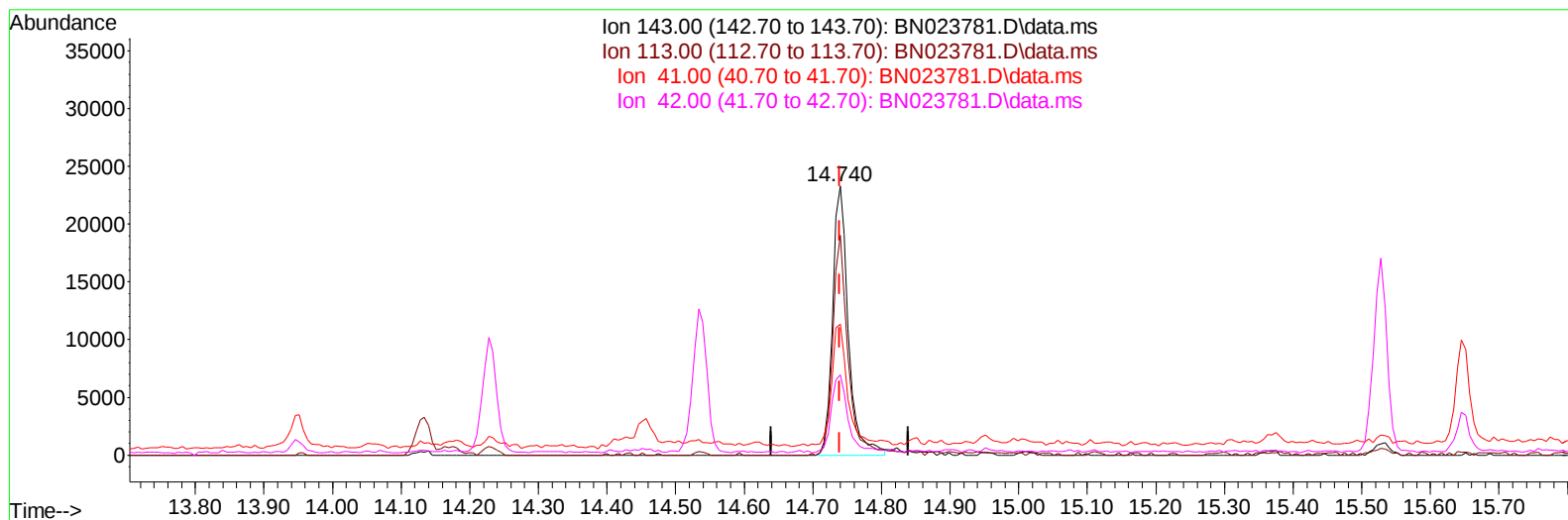
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TIC: BN023781.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (-0.000) 5.09 ng/ul m

response 37094

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.00	81.63
41.00	40.90	48.81
42.00	27.80	29.89

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	195584	20.000	ng/ul	0.00
20) Naphthalene-d8	10.699	136	810055	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	481216	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1086858	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	995047	20.000	ng/ul	0.00
88) Perylene-d12	23.898	264	1196364	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.270	96	20711	4.557	ng/uL	0.00
4) Pyridine-d5	3.699	84	204861	15.023	ng/ul	0.00
7) Phenol-d5	7.052	99	123641	7.133	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	226246	21.276	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	238451	17.962	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	211920	15.317	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128	142043	23.254	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	148691	22.995	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	254118	20.603	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	409236	21.094	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	905699	25.301	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	974112	24.560	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	37094m	5.090	ng/ul	0.00
60) Fluorene-d10	15.528	176	780758	25.525	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	124703	19.020	ng/ul	0.00
73) Anthracene-d10	17.381	188	1302093	26.772	ng/ul	0.00
81) Pyrene-d10	19.675	212	1633002	28.696	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1633194	27.605	ng/ul	0.00

Target Compounds Qvalue

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

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