

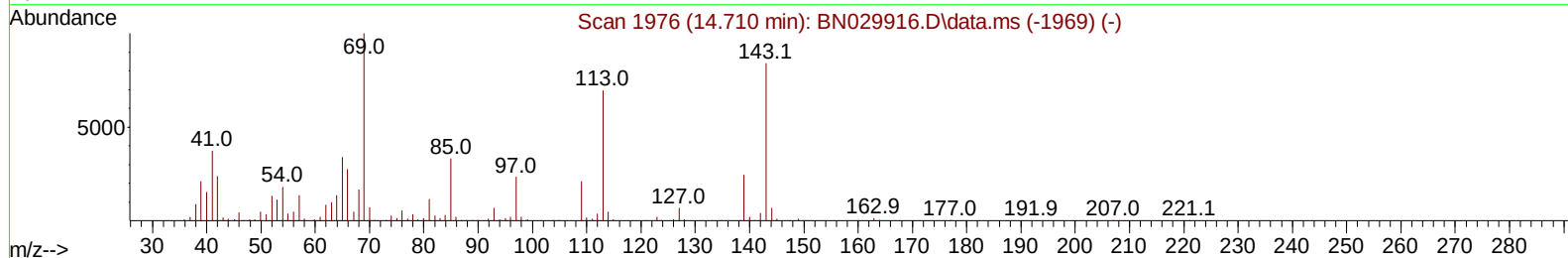
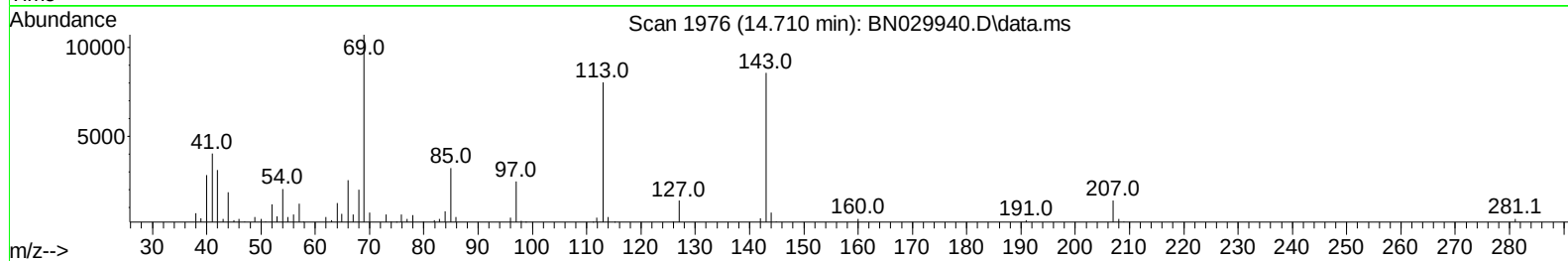
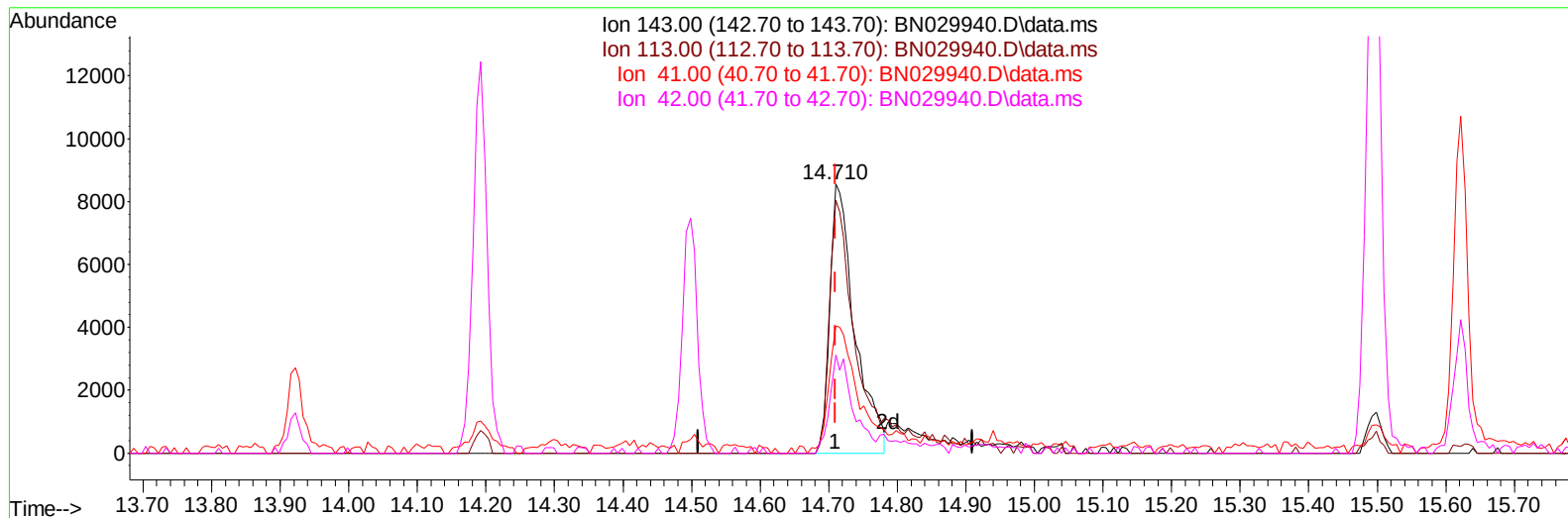
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022124\
Data File : BN029940.D
Acq On : 22 Feb 2024 22:50
Operator : MA/JU
Sample : P1494-07
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH3Z9

Manual IntegrationsAPPROVED

Quant Time: Feb 22 23:21:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 21 22:34:54 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 02/23/2024
Supervised By :mohammad ahmed 02/24/2024



TIC: BN029940.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.000) 5.48 ng/ul

response 21395

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	93.87
41.00	45.90	47.36
42.00	28.40	36.46#

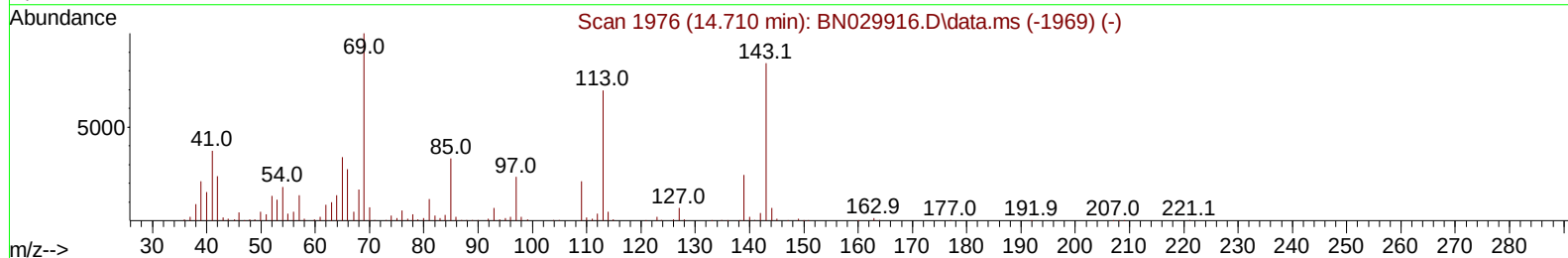
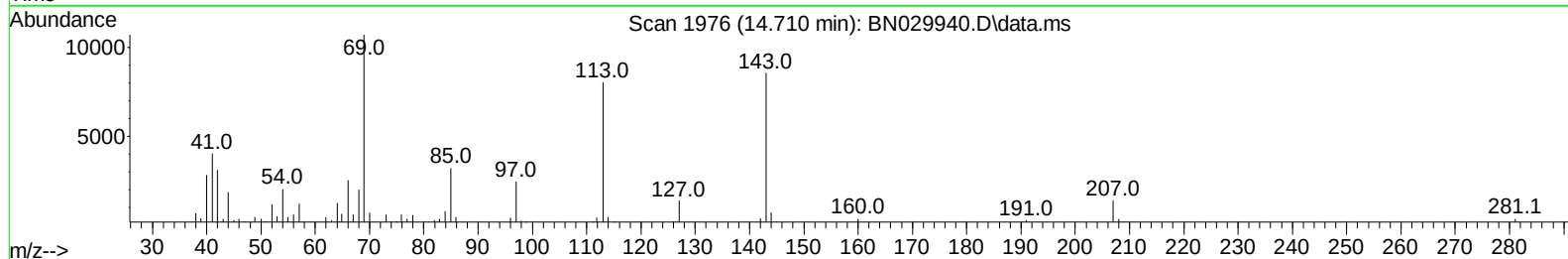
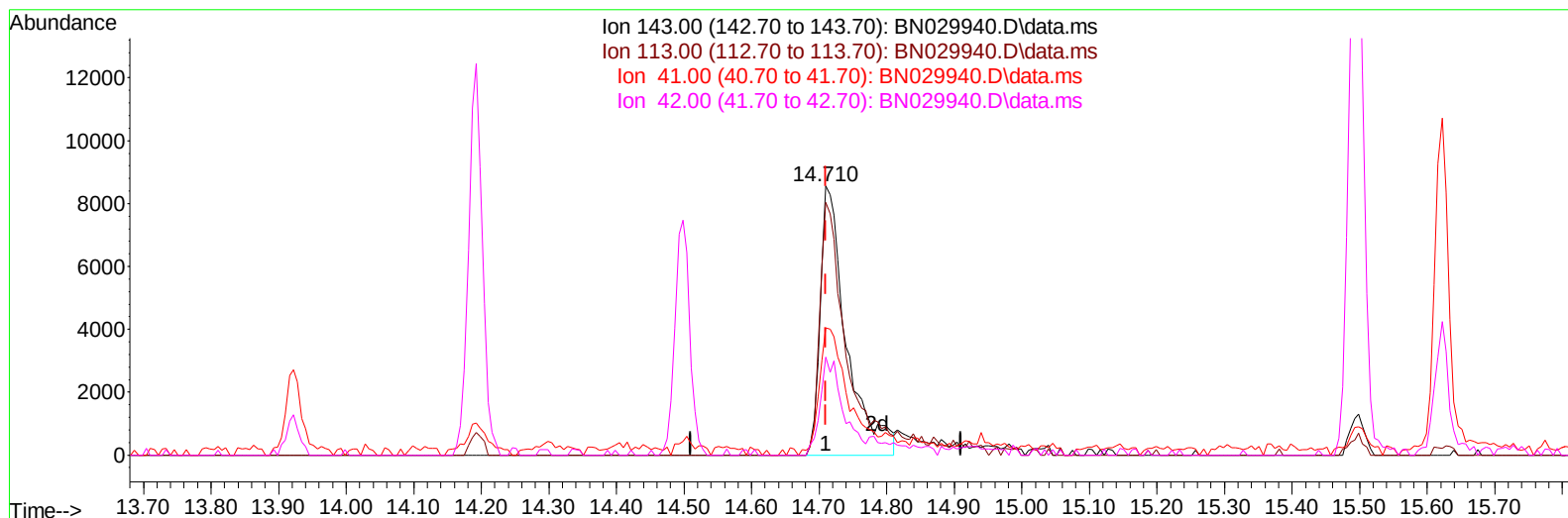
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TIC: BN029940.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.000) 5.87 ng/ul m

response 22929

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	93.87
41.00	45.90	47.36
42.00	28.40	36.46#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	129273	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	524293	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	276216	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	578159	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	508666	20.000	ng/ul	0.00
88) Perylene-d12	23.816	264	565446	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	15410	3.921	ng/uL	0.00
4) Pyridine-d5	3.734	84	73455	7.008	ng/ul	0.00
7) Phenol-d5	7.022	99	112765	9.326	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	343260	42.536	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	306737	35.007	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	197098	21.857	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	195090	47.943	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	182459	51.484	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	298823	42.718	ng/ul	0.00
31) 4-Chloroaniline-d4	10.799	131	409653	34.327	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	995319	48.400	ng/ul	0.00
49) Acenaphthylene-d8	14.193	160	1174325	48.085	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	22929m	5.875	ng/ul	0.00
60) Fluorene-d10	15.498	176	866381	49.523	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	125881	44.684	ng/ul	0.00
73) Anthracene-d10	17.351	188	1551337	57.253	ng/ul	0.00
81) Pyrene-d10	19.651	212	1792794	58.731	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.663	264	1694411	59.885	ng/ul	0.00

Target Compounds Qvalue

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

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