

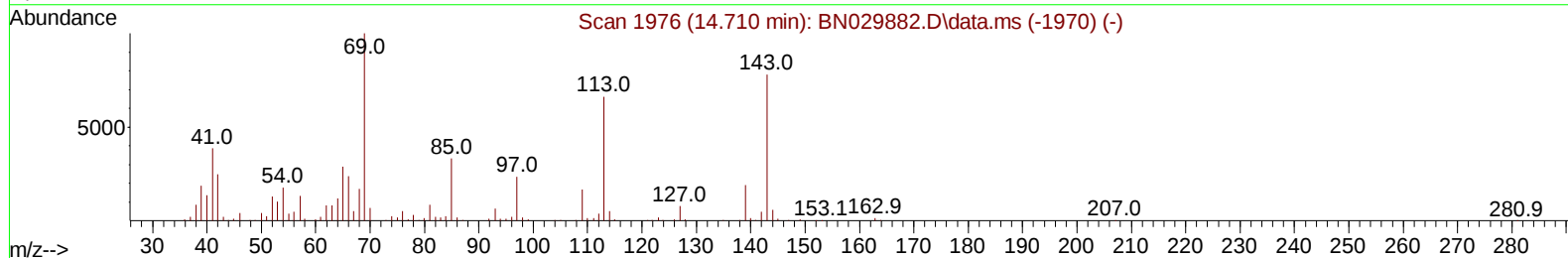
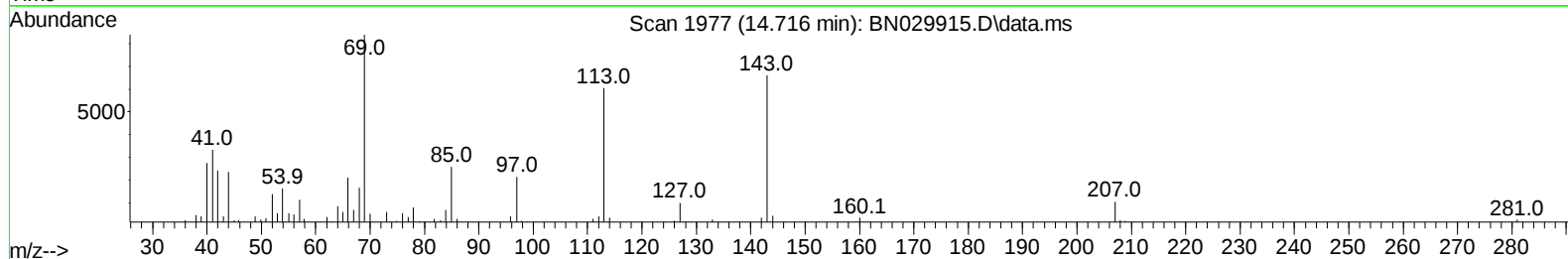
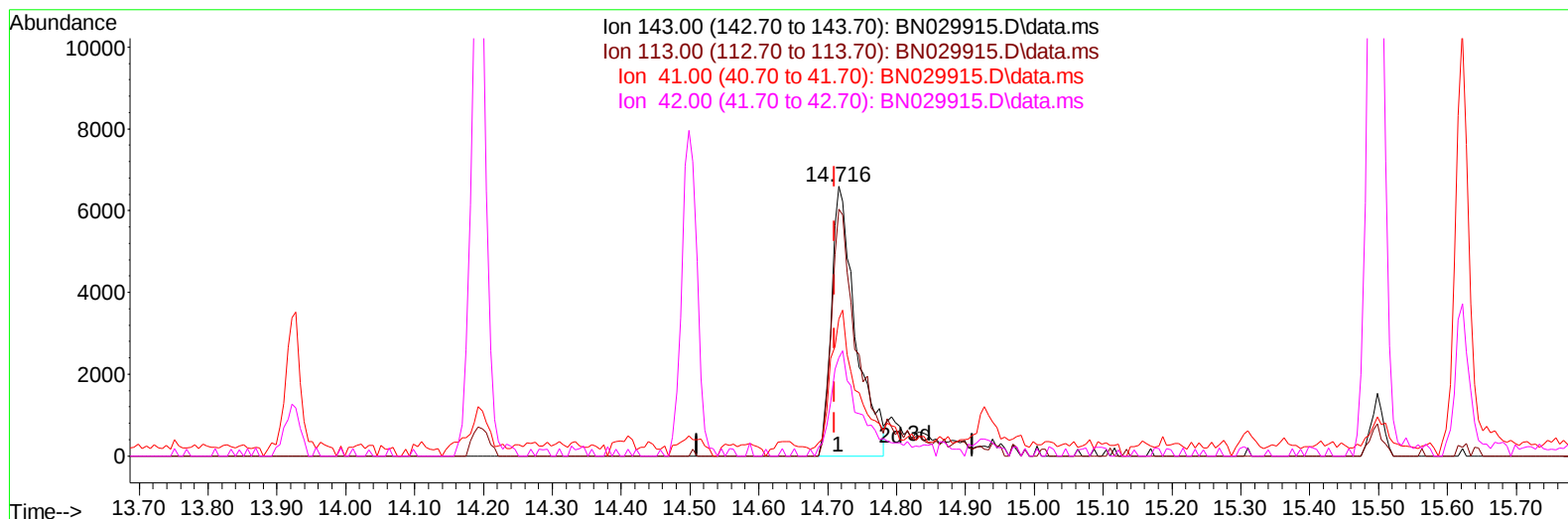
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022124\
Data File : BN029915.D
Acq On : 22 Feb 2024 06:37
Operator : MA/JU
Sample : P1493-14
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH3Y9

Manual IntegrationsAPPROVED

Quant Time: Feb 22 07:06:00 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 :Jagrut Upadhyay 02/22/2024
Supervised By :mohammad ahmed 02/24/2024



TIC: BN029915.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 3.67 ng/u1

response 16221

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	91.59
41.00	45.90	50.71
42.00	28.40	36.49#

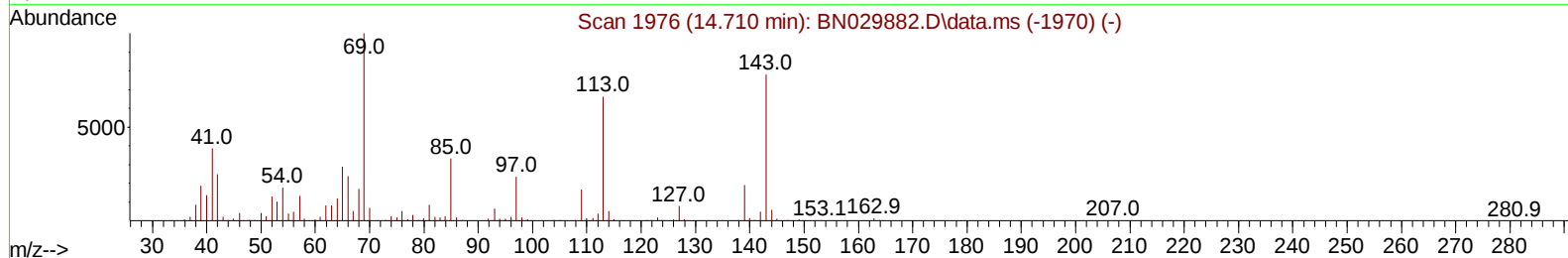
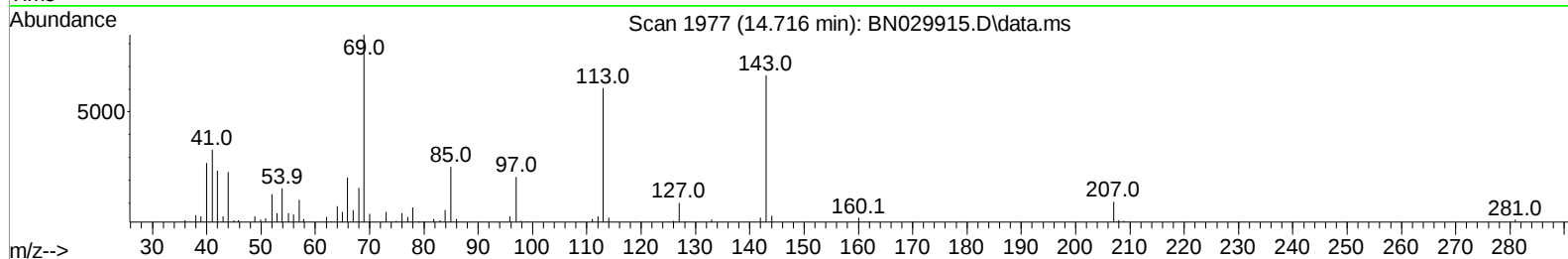
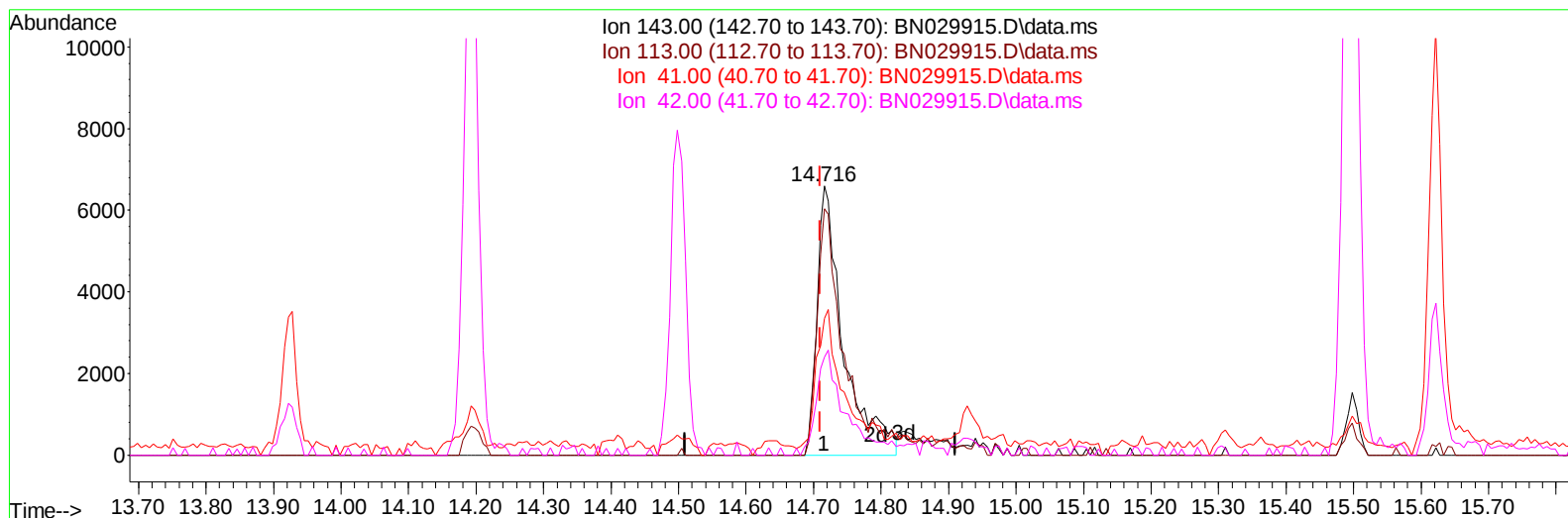
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TIC: BN029915.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.716min (+ 0.006) 4.06 ng/ul m

response 17947

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	91.59
41.00	45.90	50.71
42.00	28.40	36.49#

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/22/2024
 Supervised By :mohammad ahmed 02/24/2024

Quant Time: Feb 22 07:07:03 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	144537	20.000	ng/ul	0.00
20) Naphthalene-d8	10.652	136	585384	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	312655	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	639038	20.000	ng/ul	0.00
79) Chrysene-d12	21.457	240	568888	20.000	ng/ul	0.00
88) Perylene-d12	23.816	264	657174	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	16936	3.854	ng/uL	0.00
4) Pyridine-d5	3.734	84	92478	7.891	ng/ul	0.00
7) Phenol-d5	7.022	99	101658	7.519	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	385999	42.781	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	328890	33.571	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	191216	18.965	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	216322	47.613	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	198716	50.220	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	329838	42.231	ng/ul	0.00
31) 4-Chloroaniline-d4	10.804	131	448434	33.655	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	1118691	48.059	ng/ul	0.00
49) Acenaphthylene-d8	14.193	160	1295593	46.868	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	17947m	4.062	ng/ul	0.00
60) Fluorene-d10	15.498	176	981836	49.581	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	114169	36.666	ng/ul	0.00
73) Anthracene-d10	17.351	188	1505721	50.276	ng/ul	0.00
81) Pyrene-d10	19.651	212	1715862	50.260	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.668	264	1652296	50.246	ng/ul	0.00

Target Compounds Qvalue

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

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