

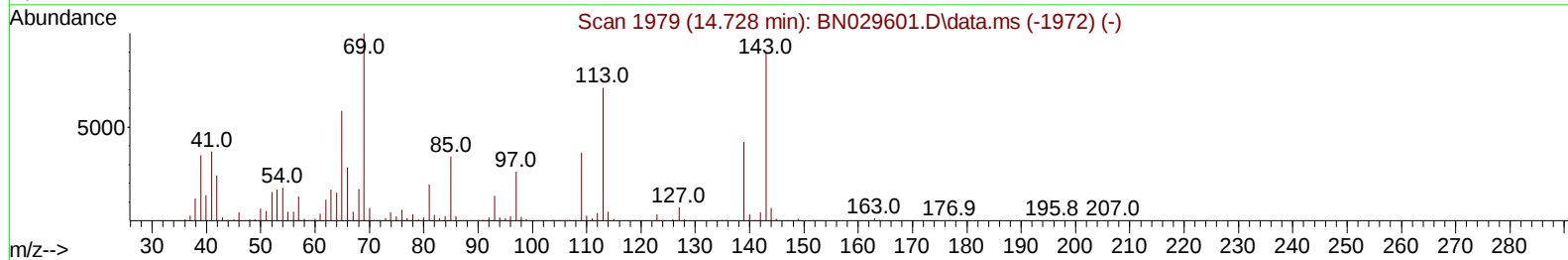
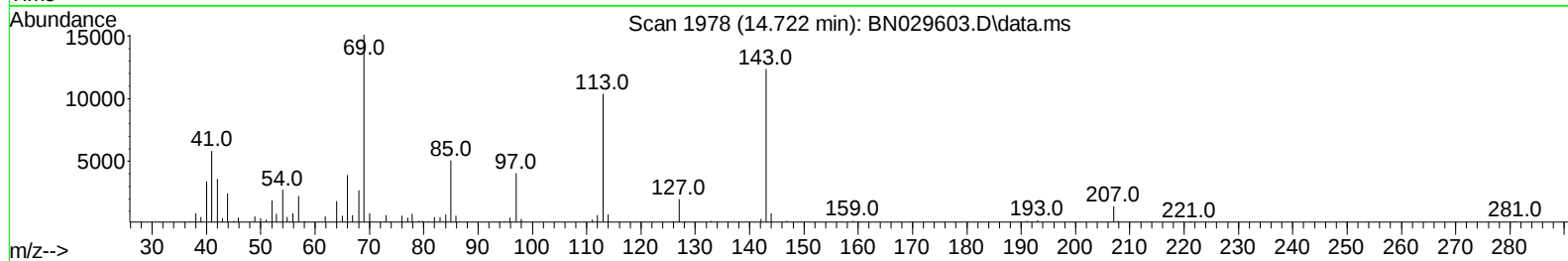
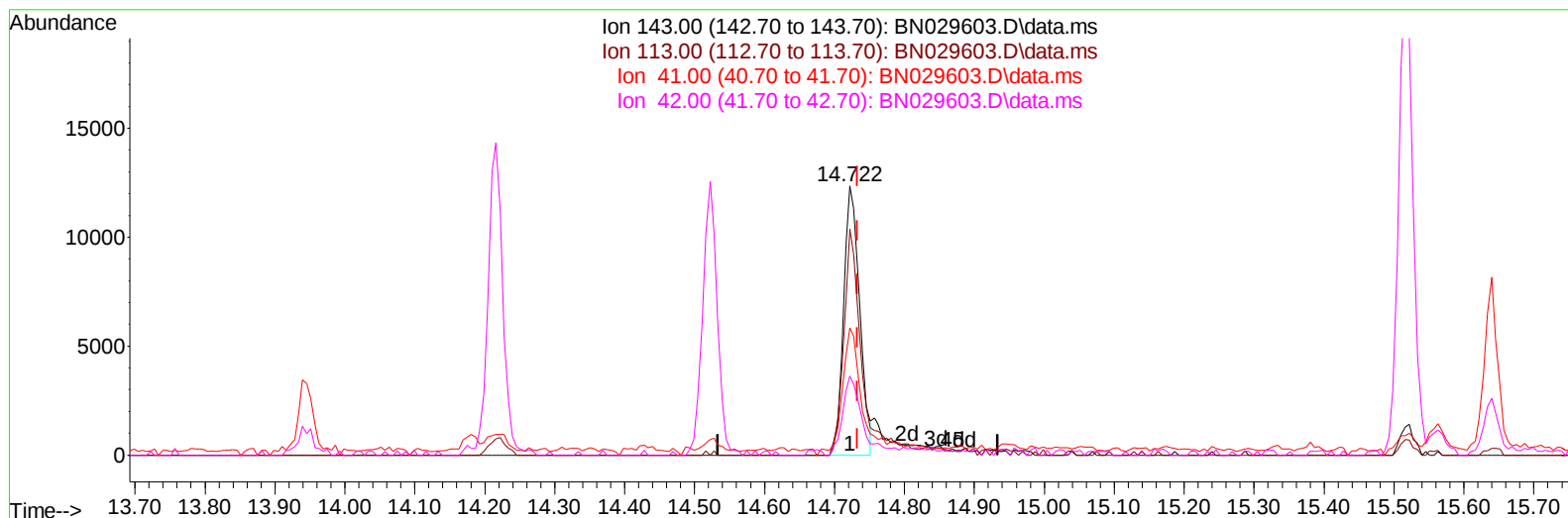
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020924\
Data File : BN029603.D
Acq On : 09 Feb 2024 10:56
Operator : MA/JU
Sample : P1300-10
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BH700

Manual IntegrationsAPPROVED

Quant Time: Feb 09 11:32:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Feb 08 00:25:29 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/12/2024
Supervised By :mohammad ahmed 02/13/2024



TIC: BN029603.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.012) 2.92 ng/u1

response 19929

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	84.08
41.00	45.90	47.36
42.00	28.40	29.28

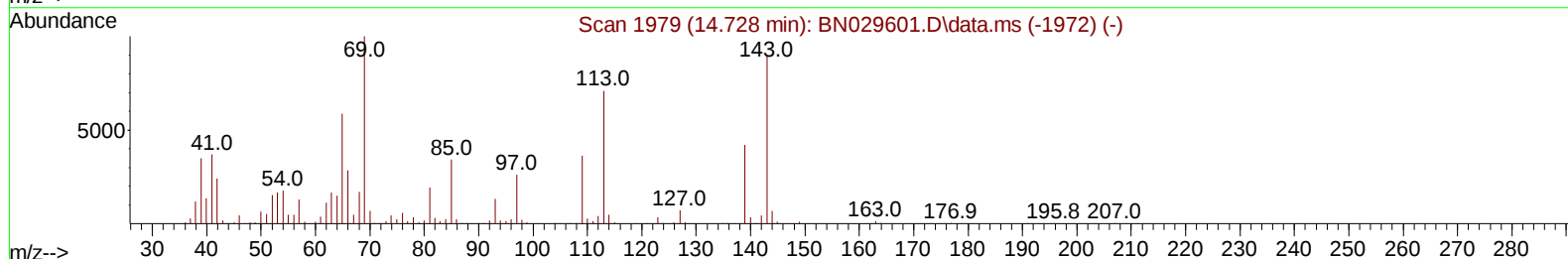
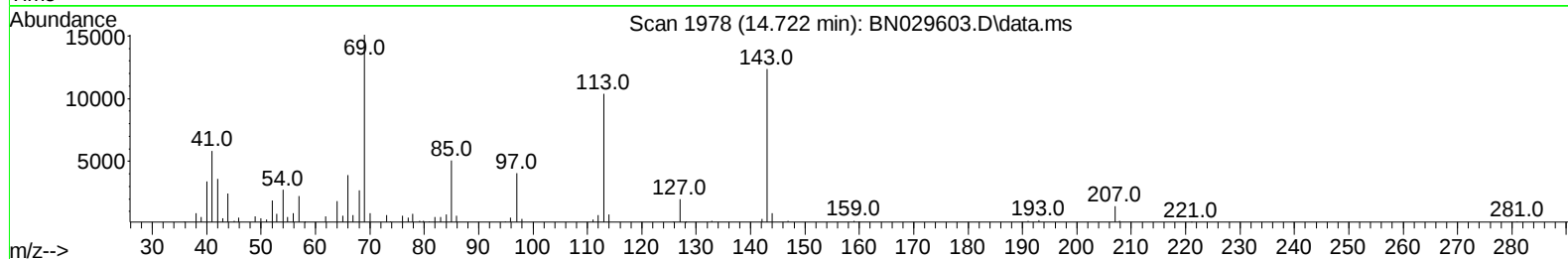
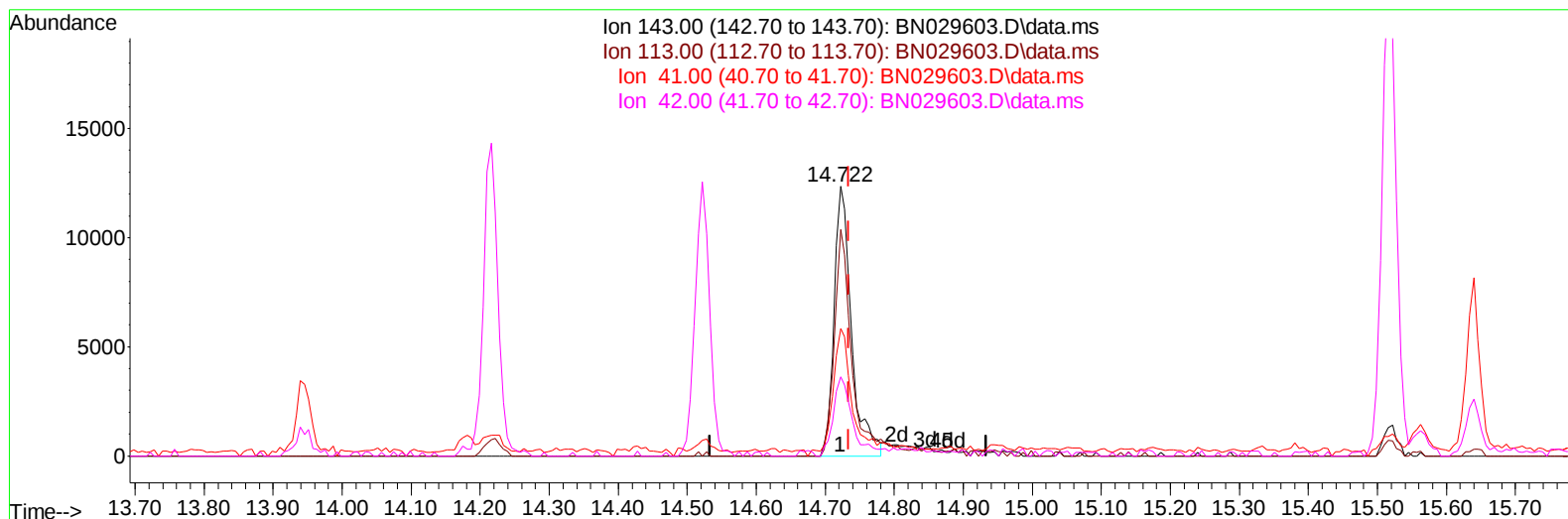
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TIC: BN029603.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.012) 3.19 ng/ul m

response 21764

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	84.08
41.00	45.90	47.36
42.00	28.40	29.28

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/12/2024
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Quant Time: Feb 09 21:59:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 08 00:25:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	242704	20.000	ng/ul	0.00
20) Naphthalene-d8	10.675	136	984663	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	483360	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	841550	20.000	ng/ul	0.00
79) Chrysene-d12	21.474	240	652097	20.000	ng/ul	0.00
88) Perylene-d12	23.845	264	737498	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	28207	3.823	ng/uL	0.00
4) Pyridine-d5	3.740	84	146754	7.458	ng/ul	0.00
7) Phenol-d5	7.046	99	160599	7.074	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	510083	33.667	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	448905	27.288	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	285770	16.879	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	271460	35.521	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	231389	34.765	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	401384	30.552	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	380136	16.961	ng/ul	-0.01
46) Dimethylphthalate-d6	13.945	166	1205345	33.494	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1486832	34.791	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	21764m	3.186	ng/ul	-0.01
60) Fluorene-d10	15.516	176	994261	32.477	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	88518	21.587	ng/ul	0.00
73) Anthracene-d10	17.375	188	1524871	38.663	ng/ul	0.00
81) Pyrene-d10	19.669	212	1577373	40.308	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1534427	41.579	ng/ul	-0.01

Target Compounds Qvalue

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

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