

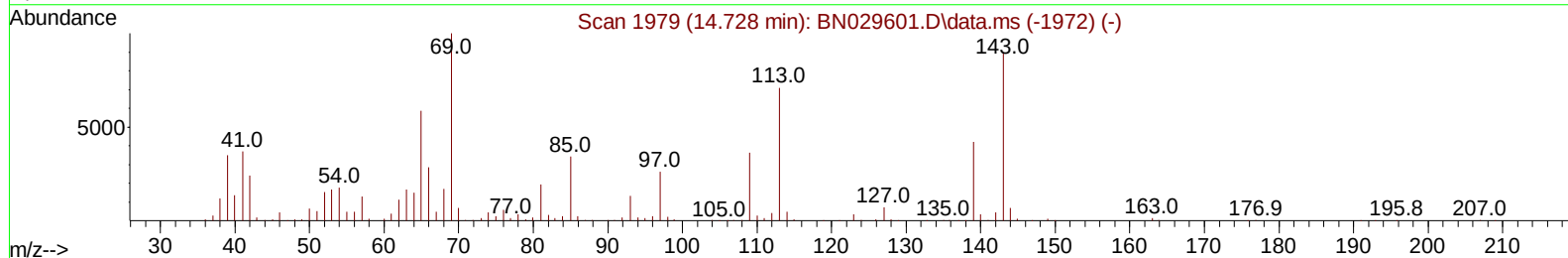
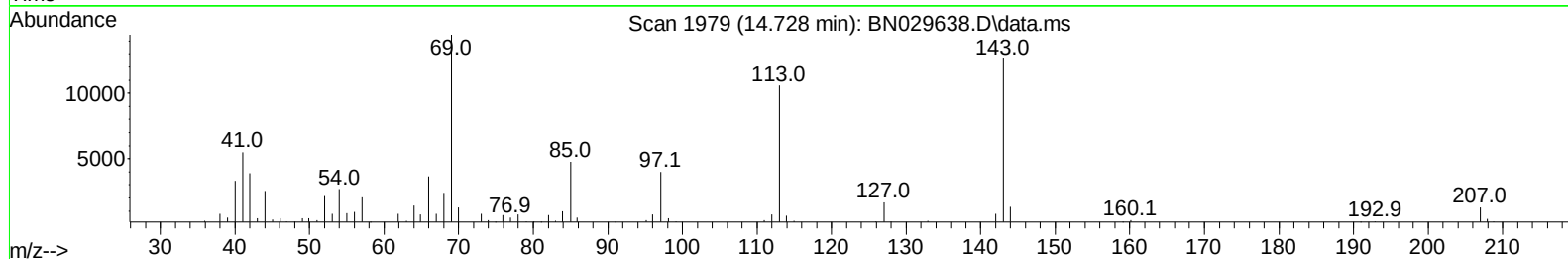
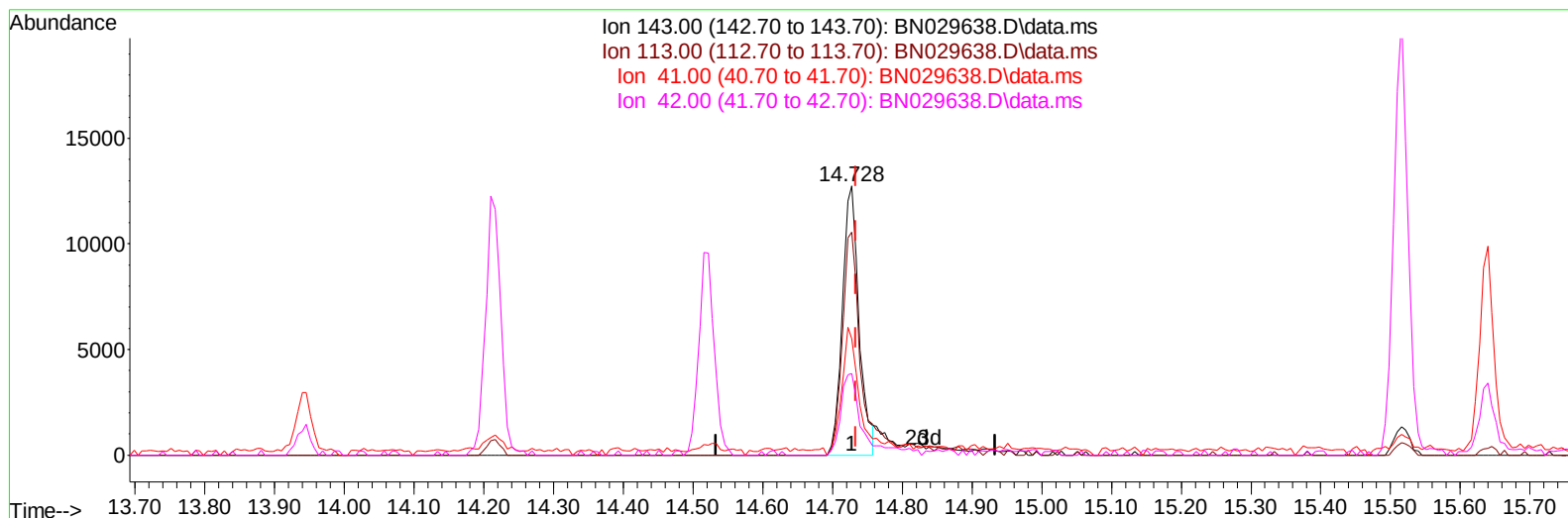
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020924\
Data File : BN029638.D
Acq On : 10 Feb 2024 08:33
Operator : MA/JU
Sample : P1399-03
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AH6

Manual IntegrationsAPPROVED

Quant Time: Feb 11 22:10:11 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: BN029638.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (-0.006) 3.79 ng/u1

response 20958

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	82.95
41.00	45.90	43.24
42.00	28.40	30.39

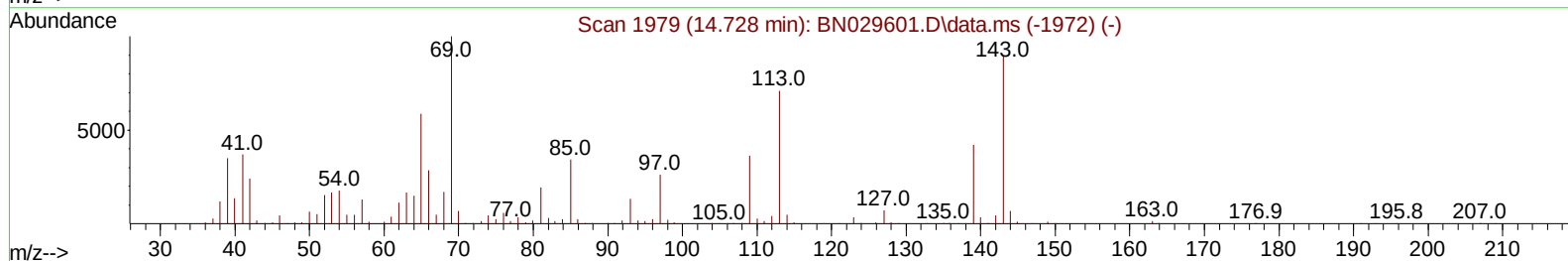
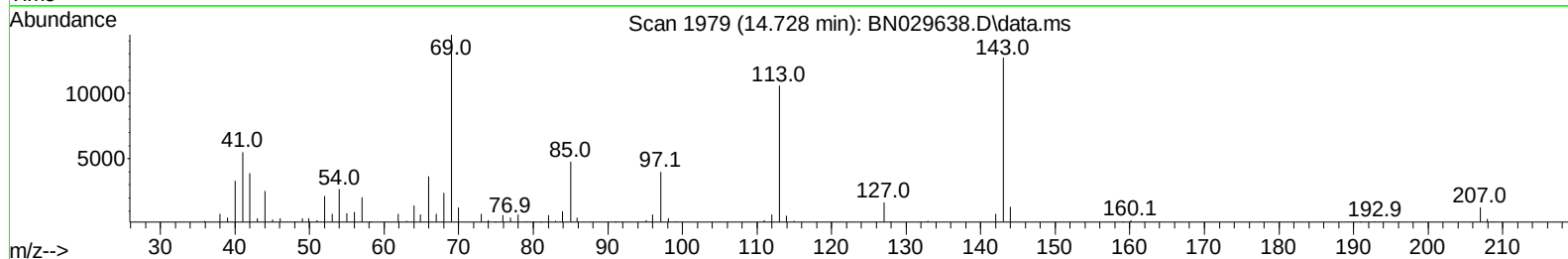
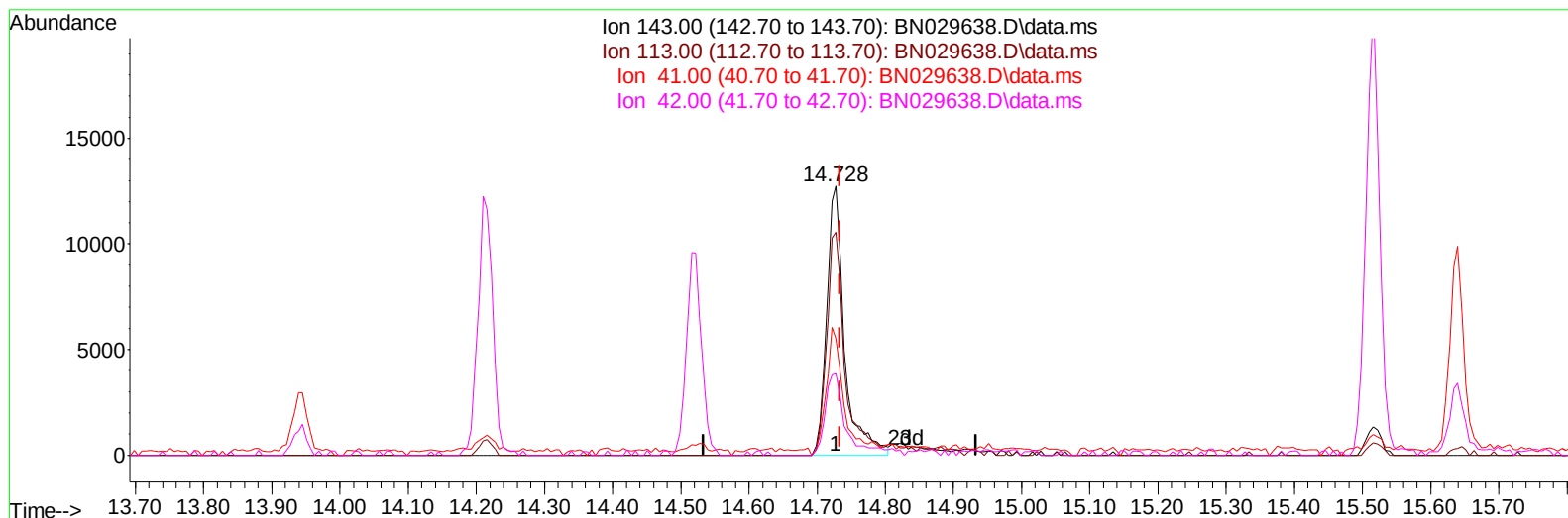
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(54) 4-Nitrophenol-d4 (S)

14.728min (-0.006) 4.17 ng/ul m

response 23092

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	82.95
41.00	45.90	43.24
42.00	28.40	30.39

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 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	7.875	152	198137	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.675	136	806007	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	391445	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	702098	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240	554959	20.000	ng/ul	-0.01
88) Perylene-d12	23.845	264	636849	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	23769	3.946	ng/uL	0.00
4) Pyridine-d5	3.746	84	71100	4.426	ng/ul	0.00
7) Phenol-d5	7.046	99	121494	6.556	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	418327	33.822	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	356192	26.522	ng/ul	-0.01
15) 4-Methylphenol-d8	8.587	113	223295	16.156	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	225857	36.104	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	198259	36.390	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.298	165	332815	30.948	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	377566	20.580	ng/ul	-0.01
46) Dimethylphthalate-d6	13.945	166	1079759	37.050	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1274576	36.827	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	23092m	4.175	ng/ul	0.00
60) Fluorene-d10	15.516	176	921825	37.181	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	116316	34.000	ng/ul	0.00
73) Anthracene-d10	17.375	188	1492904	45.371	ng/ul	0.00
81) Pyrene-d10	19.669	212	1559358	46.823	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1527323	47.928	ng/ul	-0.01

Target Compounds Qvalue

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

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