

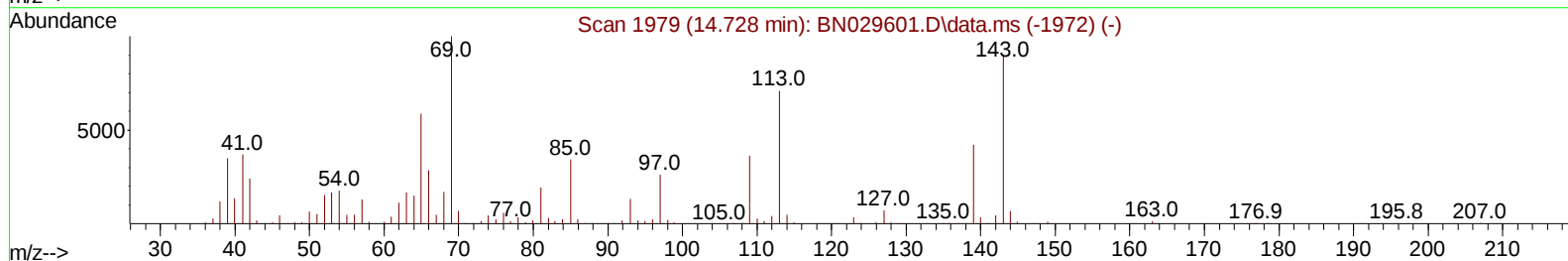
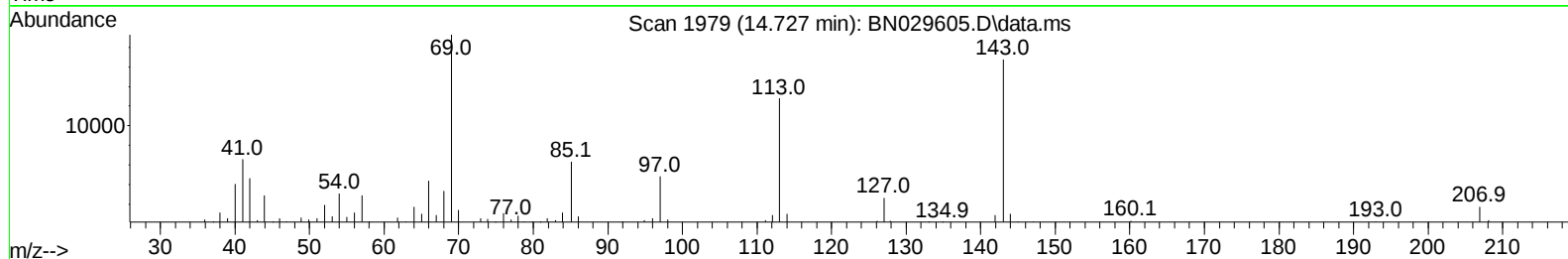
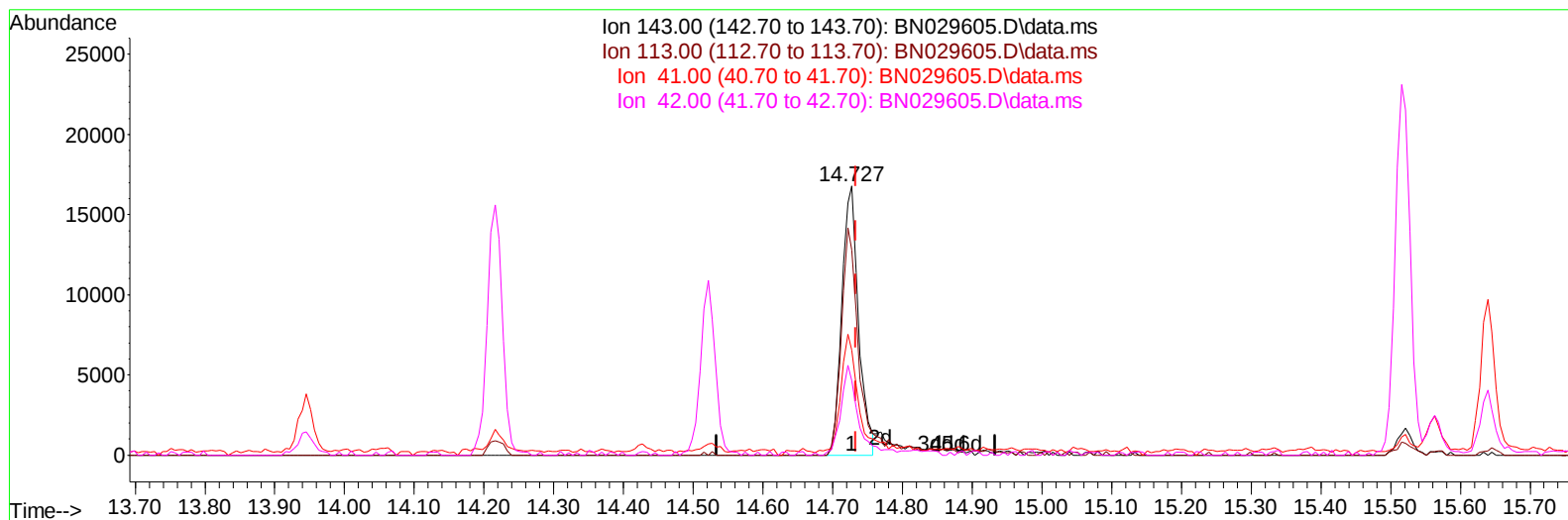
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020924\
Data File : BN029605.D
Acq On : 09 Feb 2024 12:08
Operator : MA/JU
Sample : P1348-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0R24

Manual IntegrationsAPPROVED

Quant Time: Feb 09 12:42:24 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: BN029605.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (-0.006) 4.52 ng/u1

response 27628

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	76.34
41.00	45.90	39.12
42.00	28.40	27.68

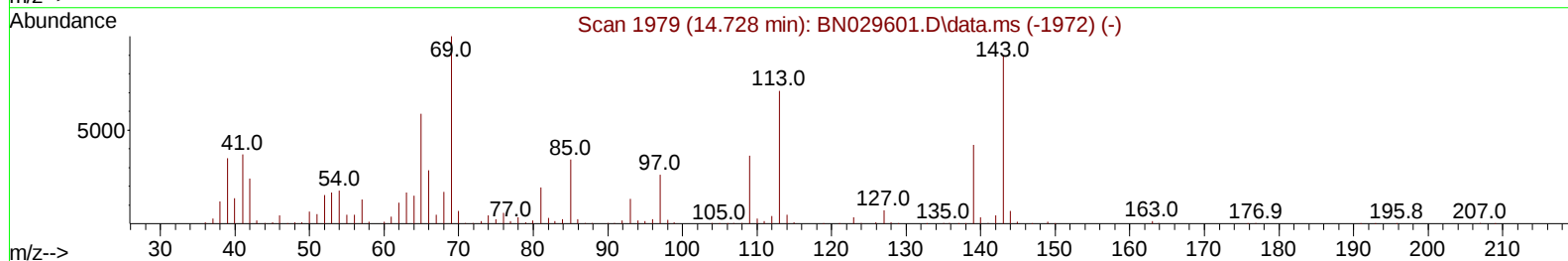
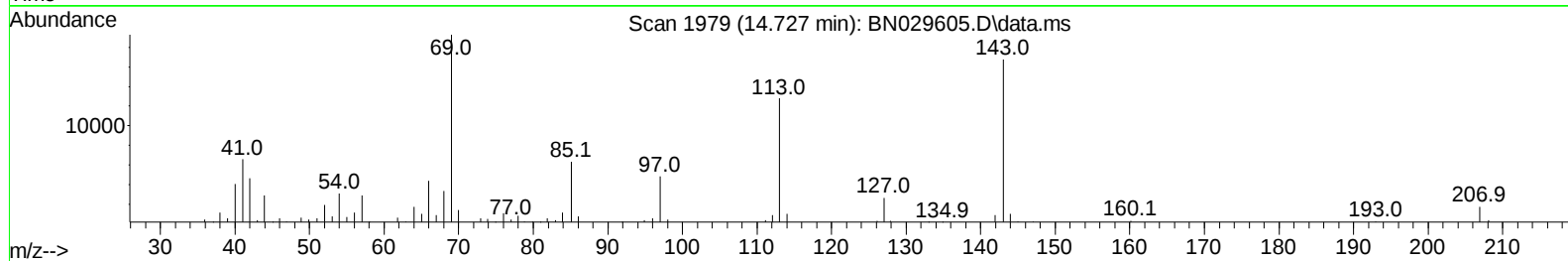
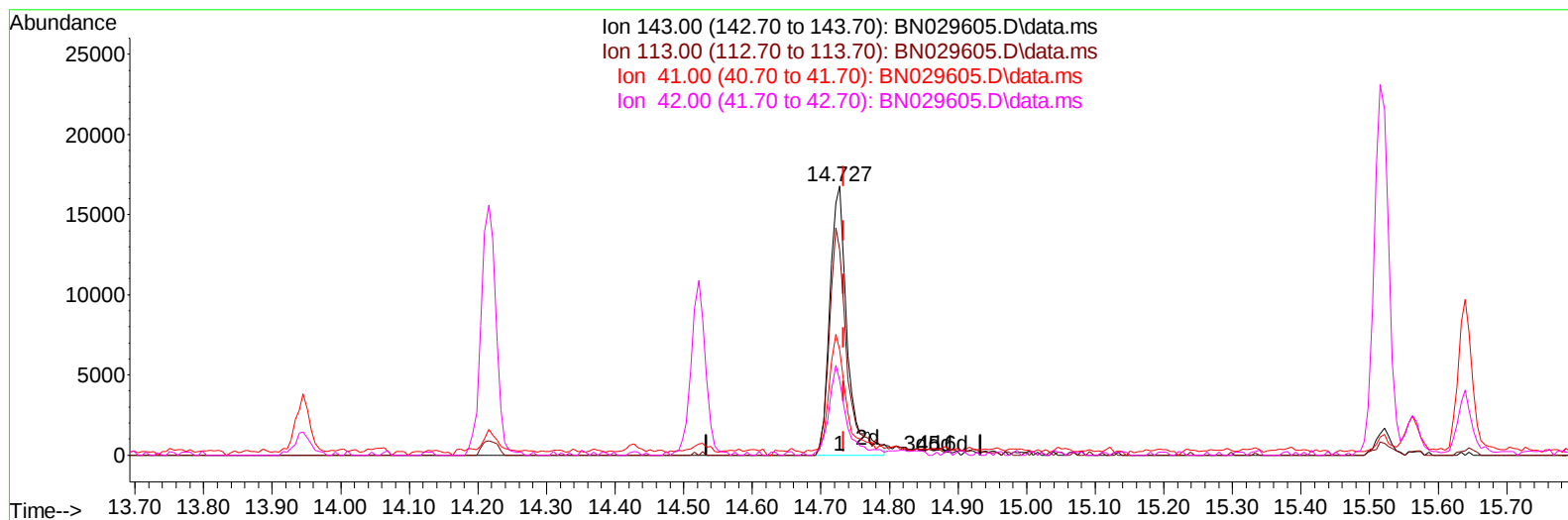
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TIC: BN029605.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (-0.006) 4.86 ng/u1 m

response 29667

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	87.40	76.34
41.00	45.90	39.12
42.00	28.40	27.68

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Manual IntegrationsAPPROVED

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Quant Time: Feb 09 22:15:49 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	222501	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	888903	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	432161	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.274	188	802664	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240	661874	20.000	ng/ul	-0.01
88) Perylene-d12	23.845	264	773273	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	29837	4.411	ng/uL	0.00
4) Pyridine-d5	3.740	84	179131	9.929	ng/ul	0.00
7) Phenol-d5	7.046	99	170496	8.192	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	543581	39.136	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	482259	31.977	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	306863	19.771	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	297735	43.156	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	262833	43.743	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	440237	37.120	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	299432	14.799	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1336759	41.547	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1614994	42.266	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	29667m	4.858	ng/ul	0.00
60) Fluorene-d10	15.522	176	1125093	41.104	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	122748	31.385	ng/ul	0.00
73) Anthracene-d10	17.374	188	1676649	44.571	ng/ul	0.00
81) Pyrene-d10	19.668	212	1802562	45.382	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1787031	46.184	ng/ul	-0.01

Target Compounds Qvalue

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

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