

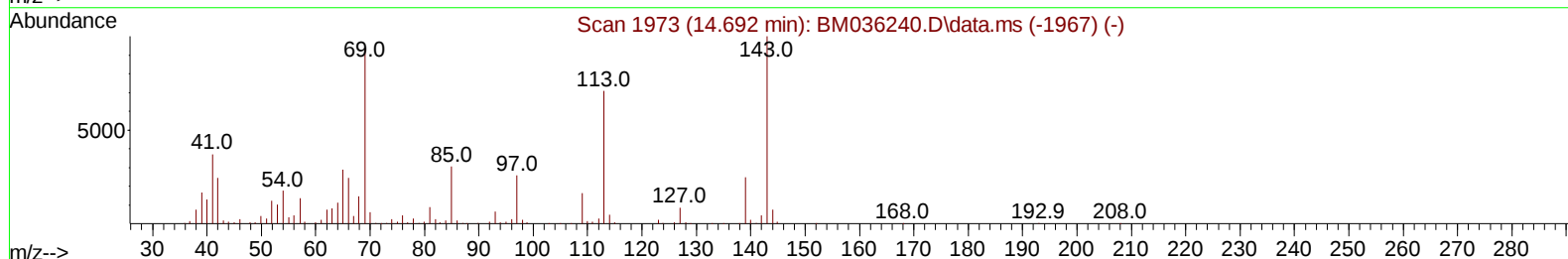
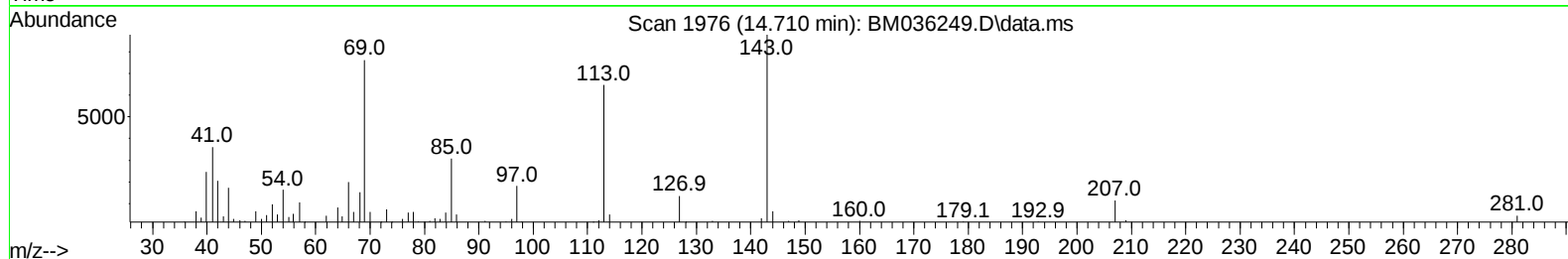
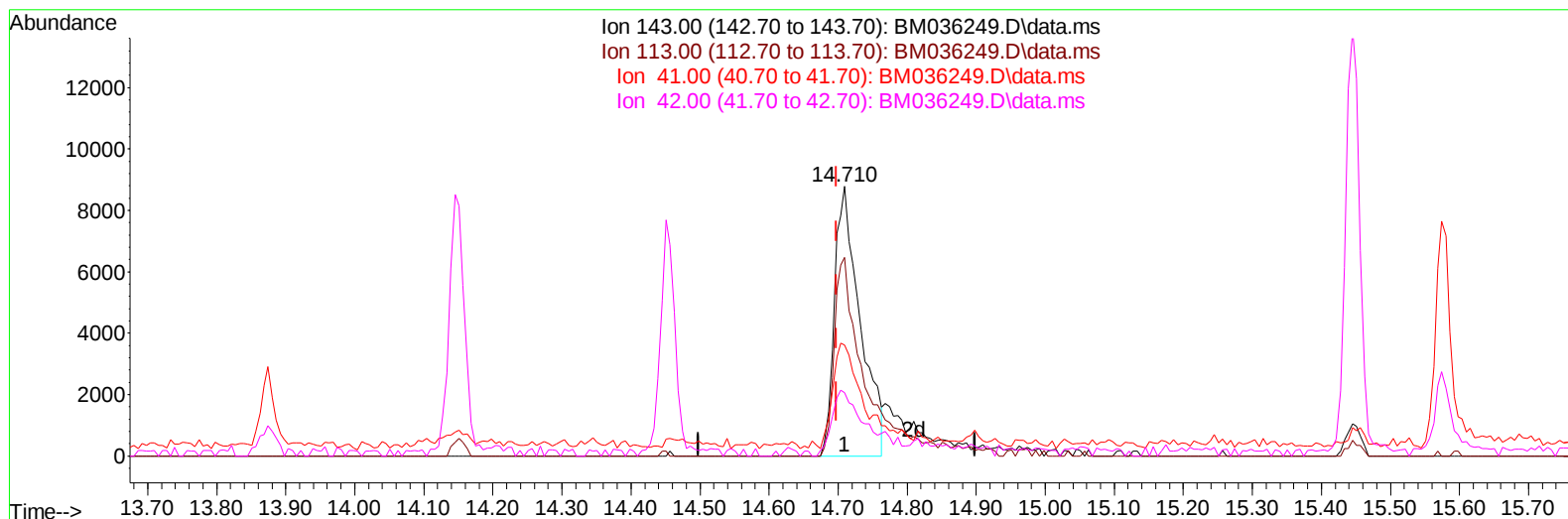
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036249.D
Acq On : 13 Aug 2022 03:11
Operator : CG/JU
Sample : N4165-03
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGBY5

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/13/2022
Supervised By :Sohil Jodhani 08/13/2022

Quant Time: Aug 13 04:06:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 12 11:23:11 2022
Response via : Initial Calibration



TIC: BM036249.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 3.98 ng/u1

response 22593

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	73.64
41.00	41.00	41.12
42.00	29.30	23.51

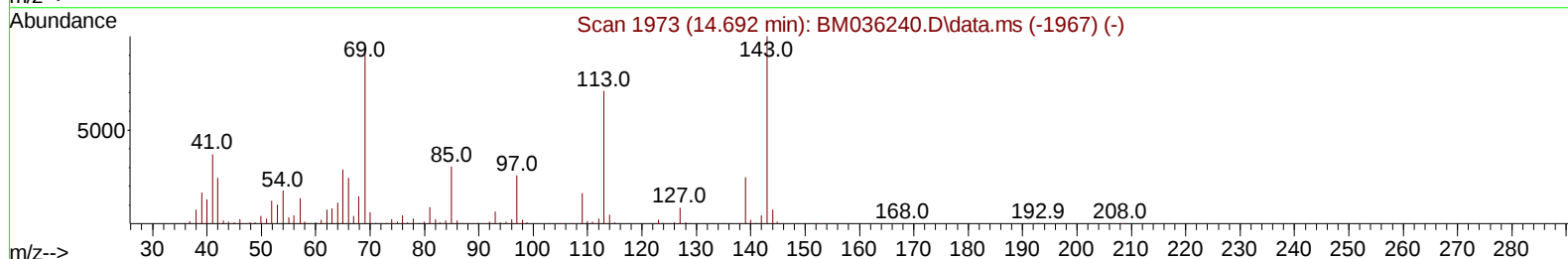
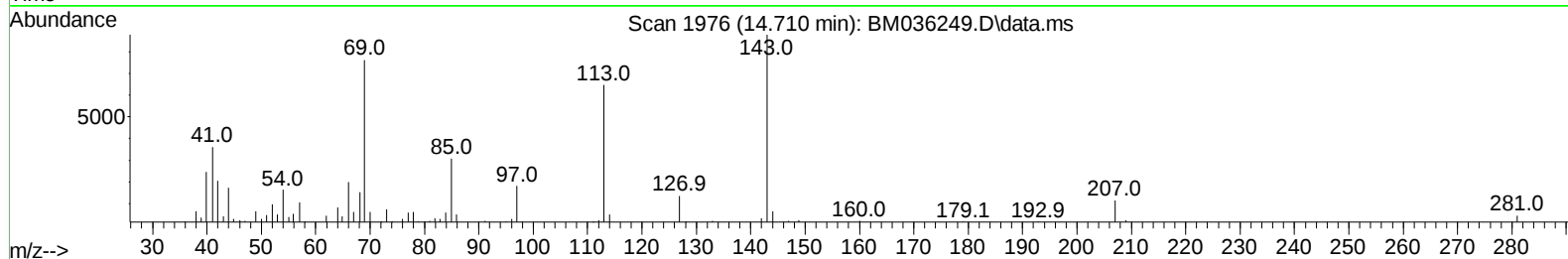
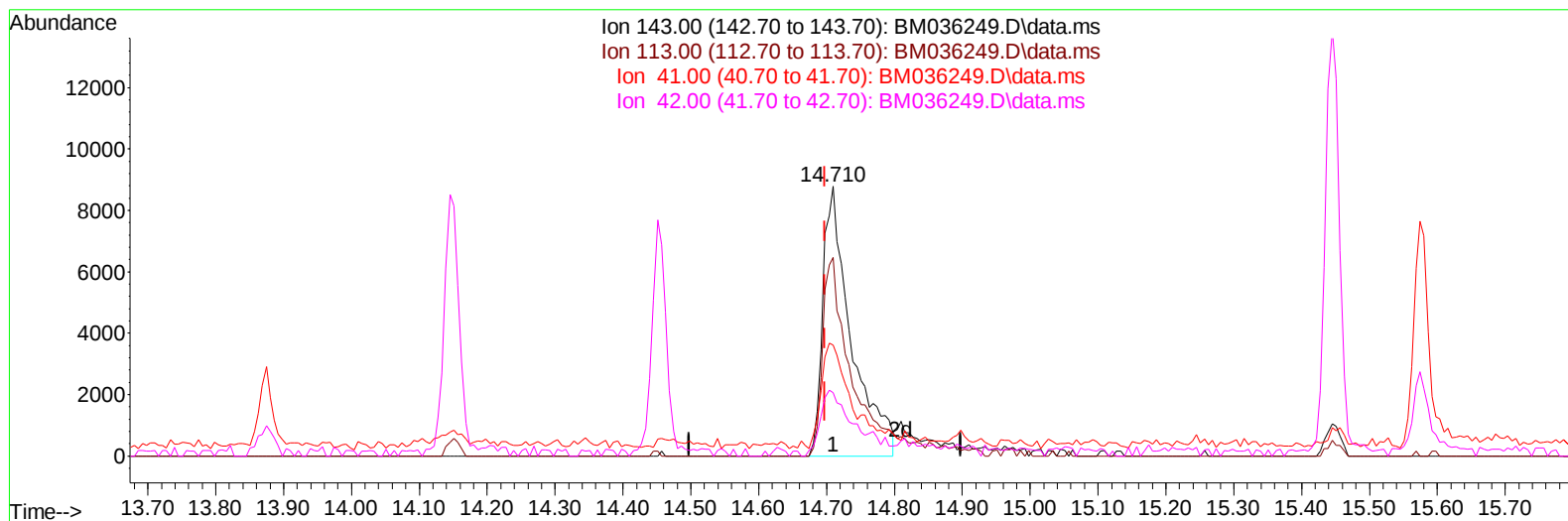
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TIC: BM036249.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 4.47 ng/ul m

response 25371

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	73.64
41.00	41.00	41.12
42.00	29.30	23.51

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	177254	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	725947	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	405681	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	827162	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	880667	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	972794	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	22231	4.719	ng/uL	0.00
4) Pyridine-d5	3.646	84	121347	9.403	ng/ul	0.00
7) Phenol-d5	7.004	99	96143	6.468	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	304028	30.681	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	304593	24.687	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113	174461	14.488	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	190418	33.102	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	184027	32.836	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	285921	28.207	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	448884	26.159	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	960837	34.770	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1159432	32.884	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	25371m	4.471	ng/ul	0.01
60) Fluorene-d10	15.445	176	894056	35.803	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	123256	27.721	ng/ul	0.00
73) Anthracene-d10	17.298	188	1465890	39.199	ng/ul	0.00
81) Pyrene-d10	19.592	212	1738099	35.611	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.568	264	1924049	38.958	ng/ul	0.00

Target Compounds Qvalue

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

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