

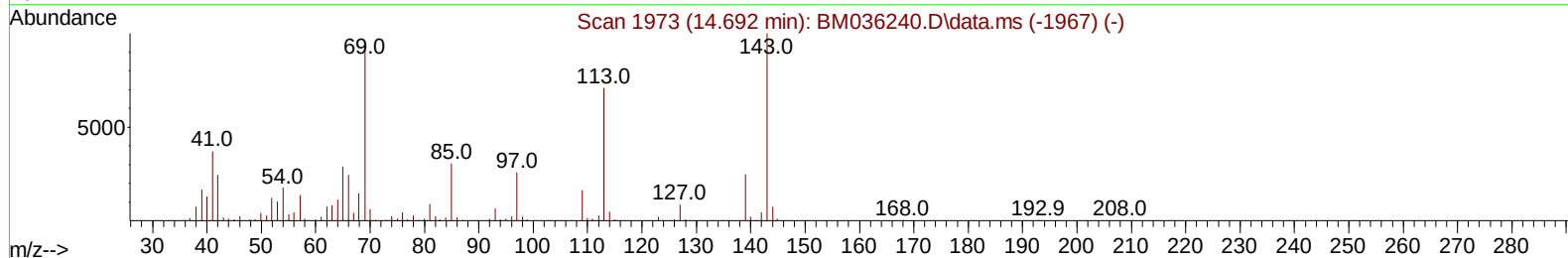
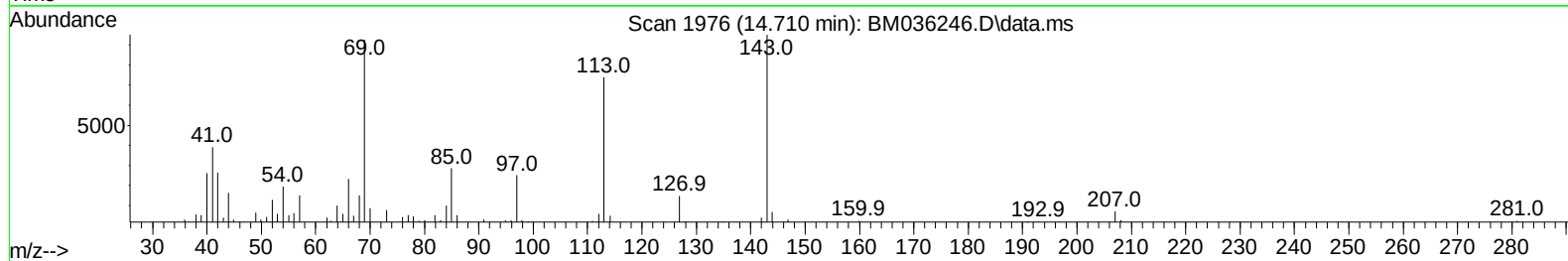
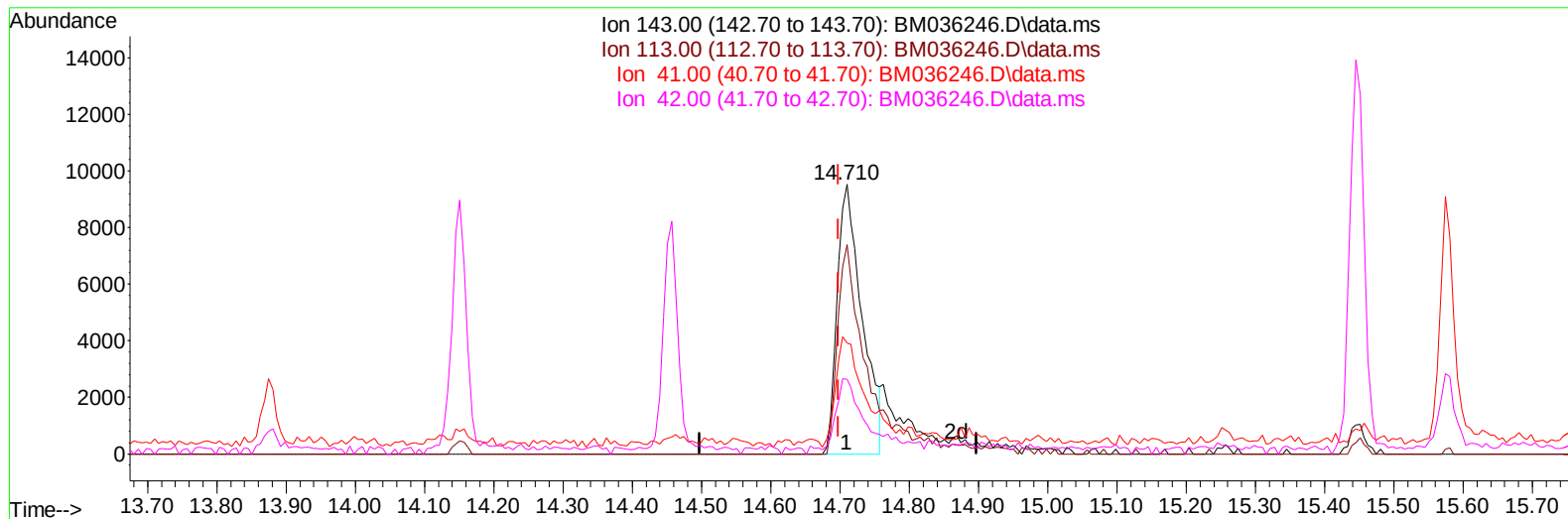
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036246.D
Acq On : 13 Aug 2022 01:24
Operator : CG/JU
Sample : N4069-17
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBJM1

Manual IntegrationsAPPROVED

Quant Time: Aug 13 01:58:52 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

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



TIC: BM036246.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 3.76 ng/ul

response 23695

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	77.79#
41.00	41.00	41.11
42.00	29.30	27.73

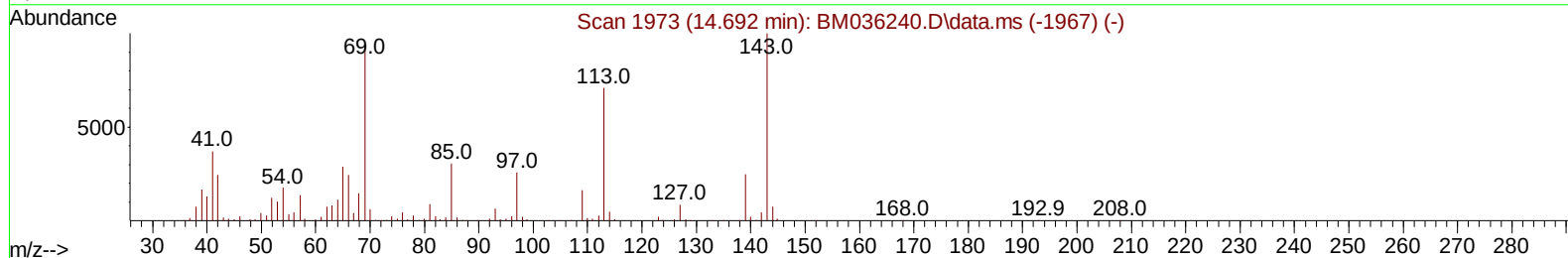
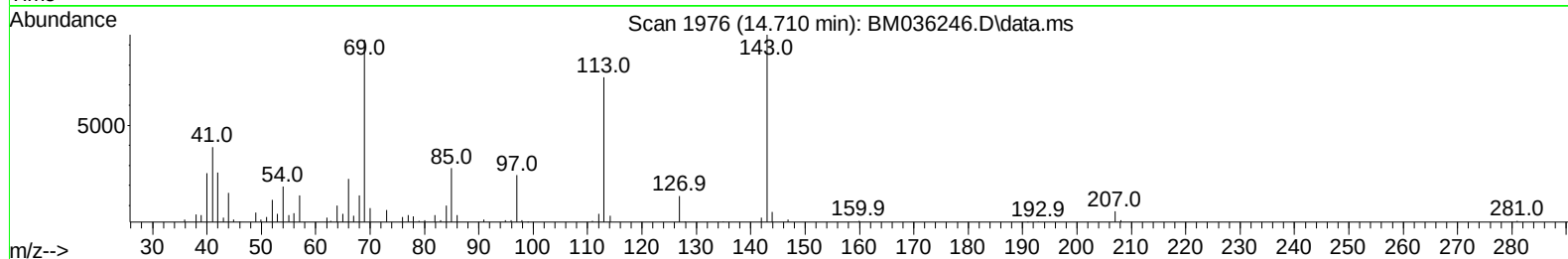
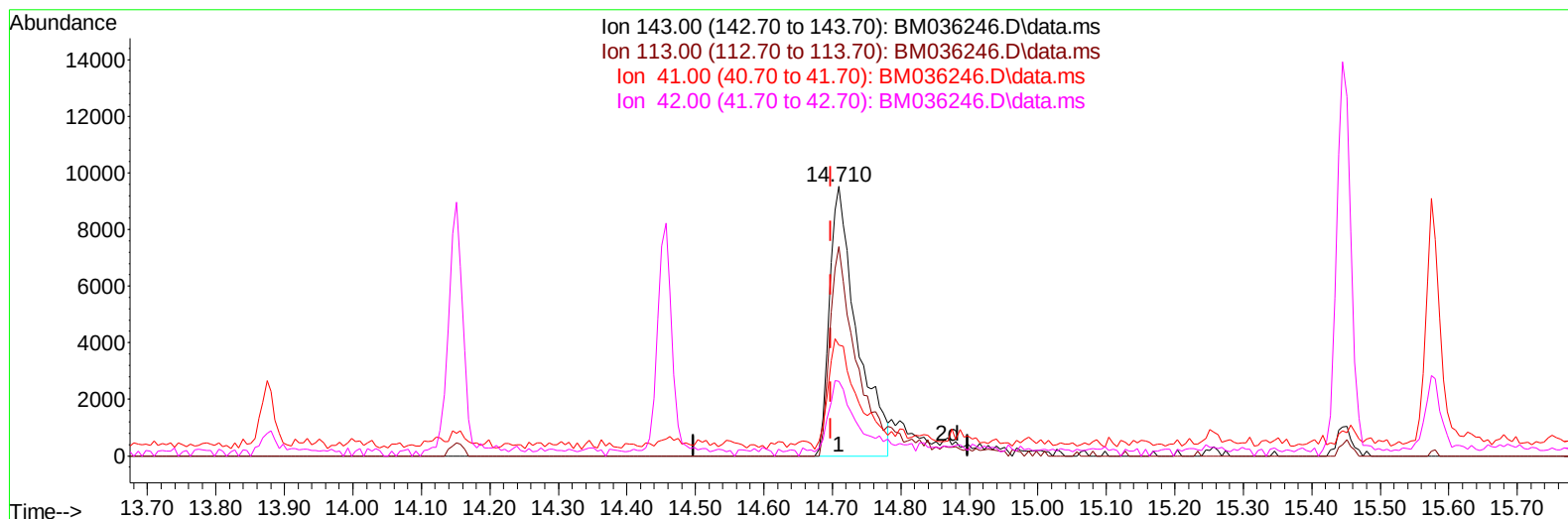
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(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 4.14 ng/ul m

response 26123

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	77.79#
41.00	41.00	41.11
42.00	29.30	27.73

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	200262	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	821517	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	451061	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	889111	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	920386	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1039716	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	23047	4.331	ng/uL	0.00
4) Pyridine-d5	3.652	84	117002	8.025	ng/ul	0.00
7) Phenol-d5	7.004	99	98639	5.874	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	302547	27.024	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	311849	22.371	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	180658	13.279	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	193459	29.718	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	184274	29.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	292881	25.533	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	437434	22.526	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	977295	31.808	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1199474	30.597	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	26123m	4.140	ng/ul	0.01
60) Fluorene-d10	15.451	176	886422	31.926	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	137382	28.746	ng/ul	0.00
73) Anthracene-d10	17.298	188	1488161	37.021	ng/ul	0.00
81) Pyrene-d10	19.592	212	1775437	34.806	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.568	264	1989443	37.690	ng/ul	0.00

Target Compounds Qvalue

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

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