

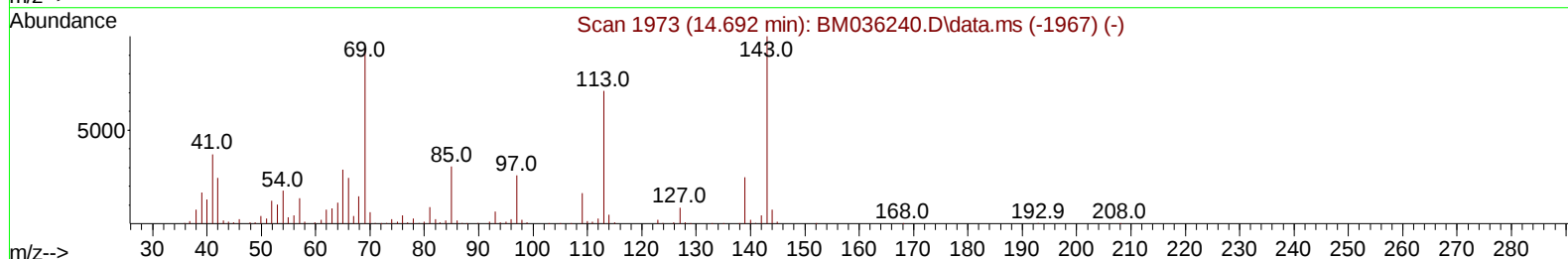
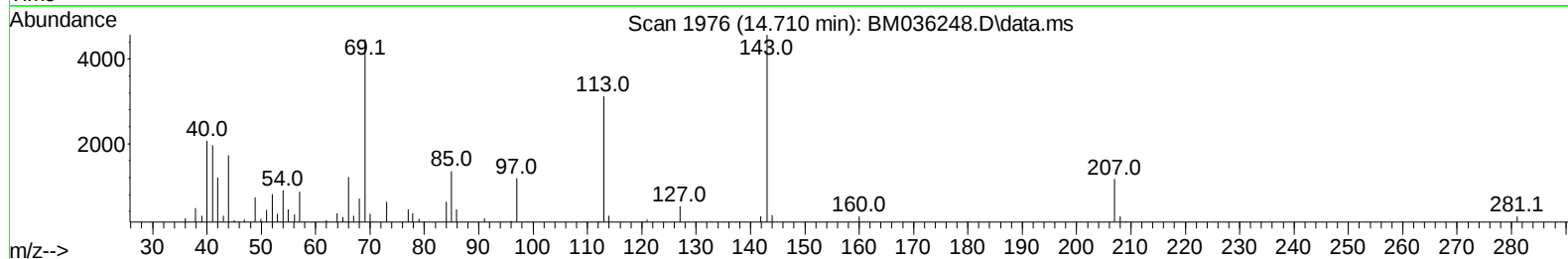
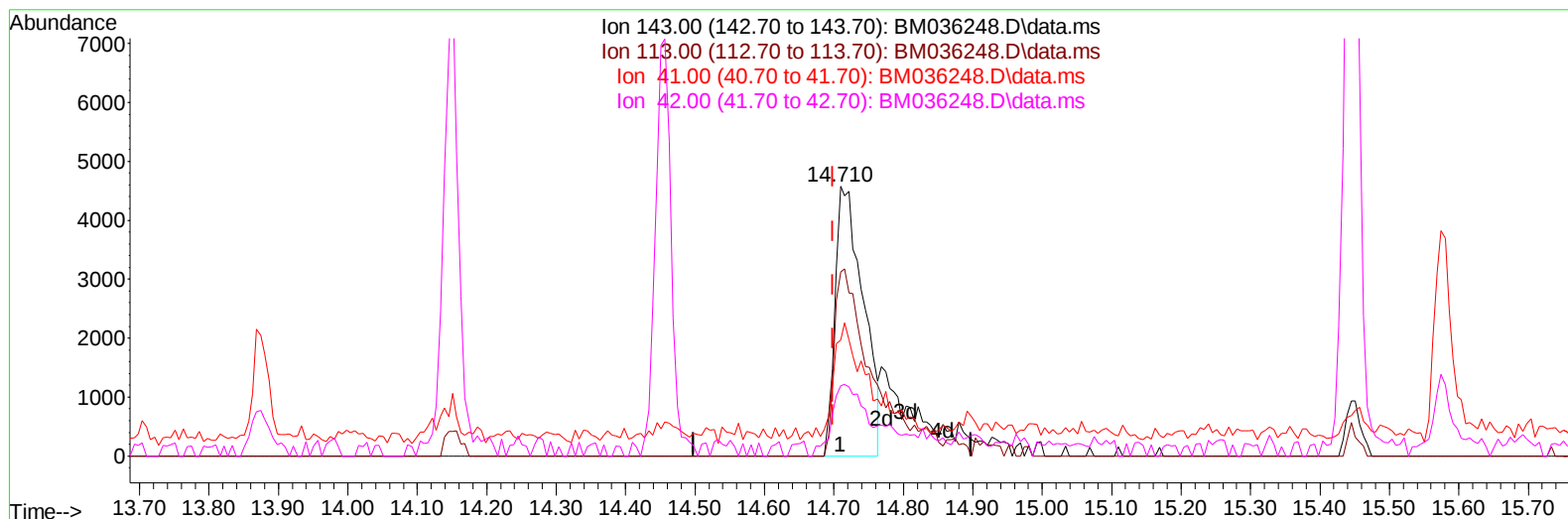
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036248.D
Acq On : 13 Aug 2022 02:36
Operator : CG/JU
Sample : N4165-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGC17

Manual IntegrationsAPPROVED

Quant Time: Aug 13 03:27:29 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: BM036248.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 2.33 ng/u1

response 12754

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	68.34
41.00	41.00	43.01
42.00	29.30	26.09

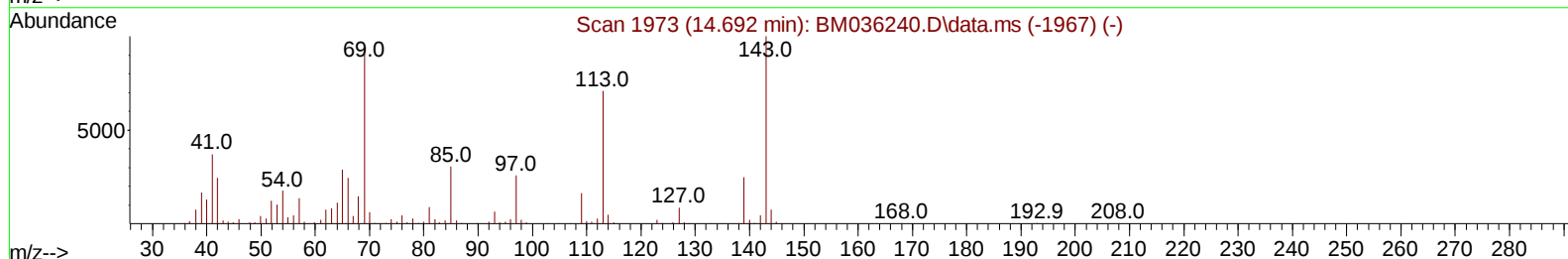
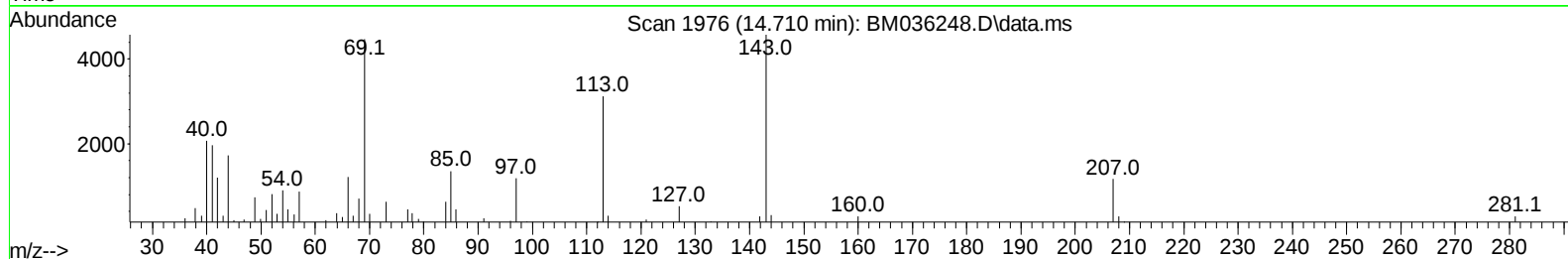
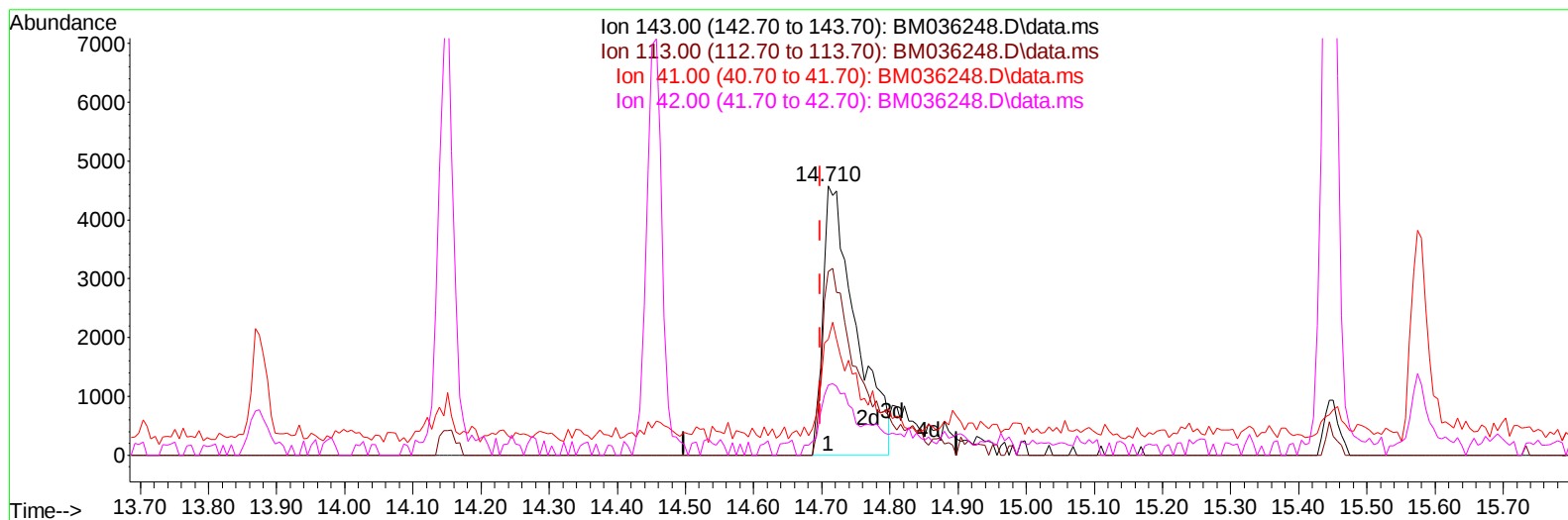
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TIC: BM036248.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.710min (+ 0.012) 2.77 ng/ul m

response 15155

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	68.34
41.00	41.00	43.01
42.00	29.30	26.09

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	170744	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	703746	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	391258	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	805630	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	876338	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	967787	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	18442	4.064	ng/uL	0.00
4) Pyridine-d5	3.658	84	51680	4.157	ng/ul	0.01
7) Phenol-d5	7.004	99	63372	4.426	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	275115	28.822	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	156672	13.182	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113	93817	8.088	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	171795	30.807	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	99299	18.277	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	150003	15.265	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	386481	23.233	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	796123	29.871	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1048900	30.846	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	15155m	2.769	ng/ul	0.01
60) Fluorene-d10	15.445	176	822289	34.143	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	62899	14.525	ng/ul	0.00
73) Anthracene-d10	17.298	188	1444068	39.647	ng/ul	0.00
81) Pyrene-d10	19.592	212	1703553	35.075	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.568	264	1938508	39.454	ng/ul	0.00

Target Compounds Qvalue

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

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