

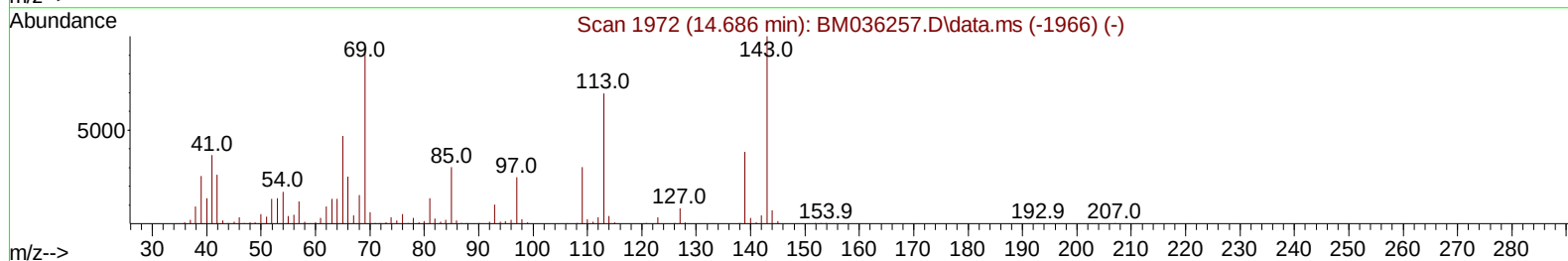
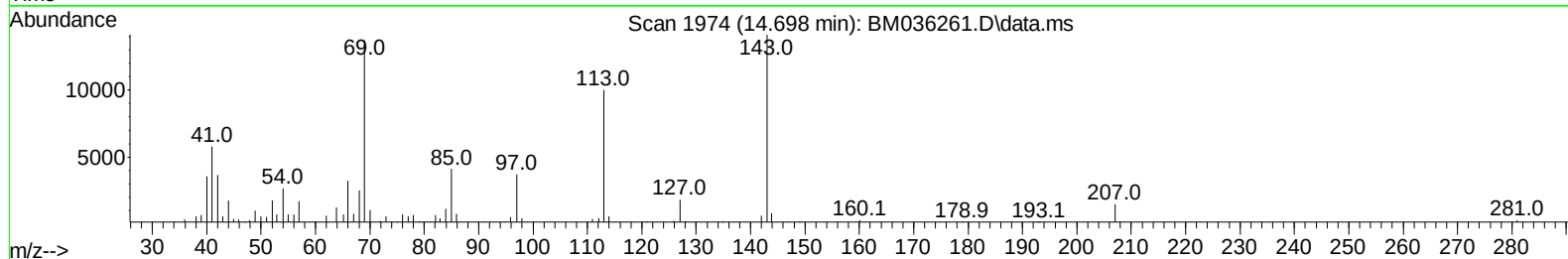
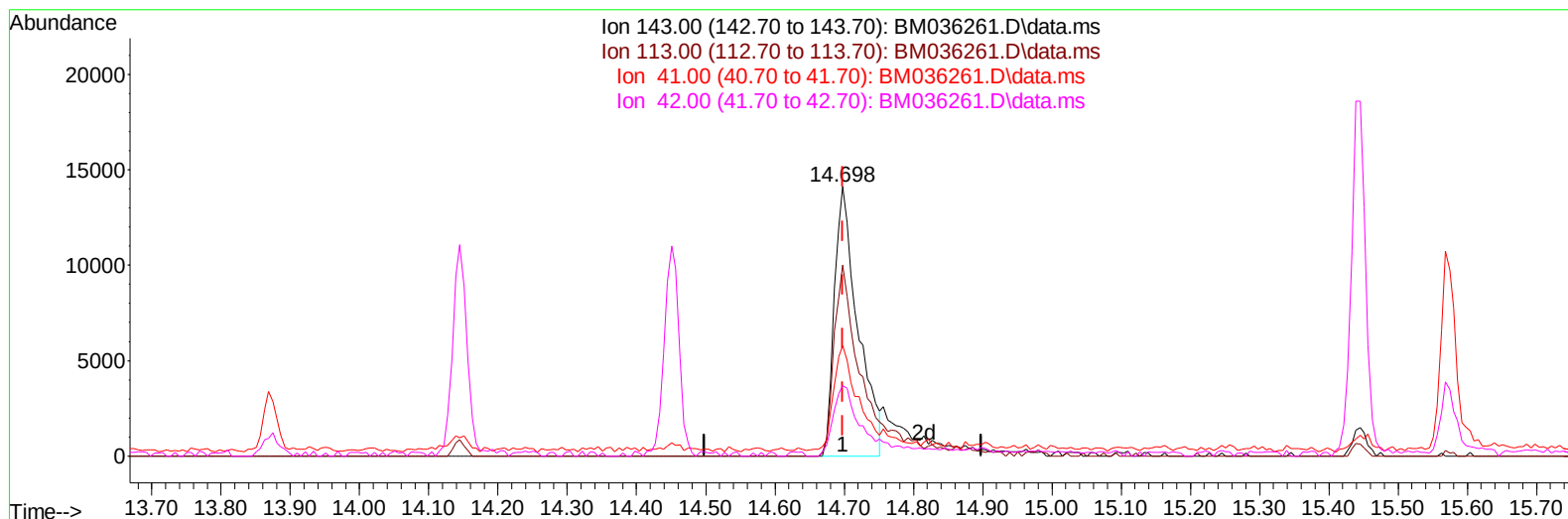
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036261.D
Acq On : 13 Aug 2022 10:57
Operator : CG/JU
Sample : N4165-08
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGBY1

Manual IntegrationsAPPROVED

Quant Time: Aug 14 23:01:01 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 08/16/2022
Supervised By :Sohil Jodhani 08/16/2022



TIC: BM036261.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.698min (-0.000) 4.06 ng/u1

response 33184

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	70.78
41.00	41.00	40.85
42.00	29.30	26.04

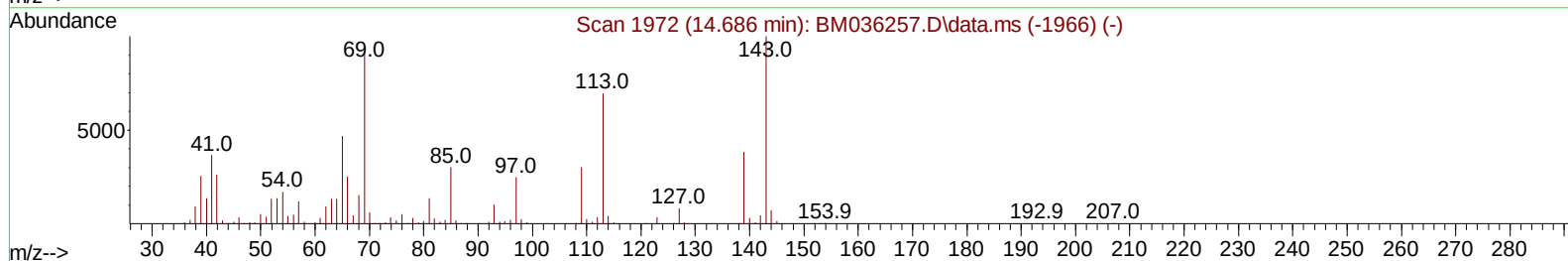
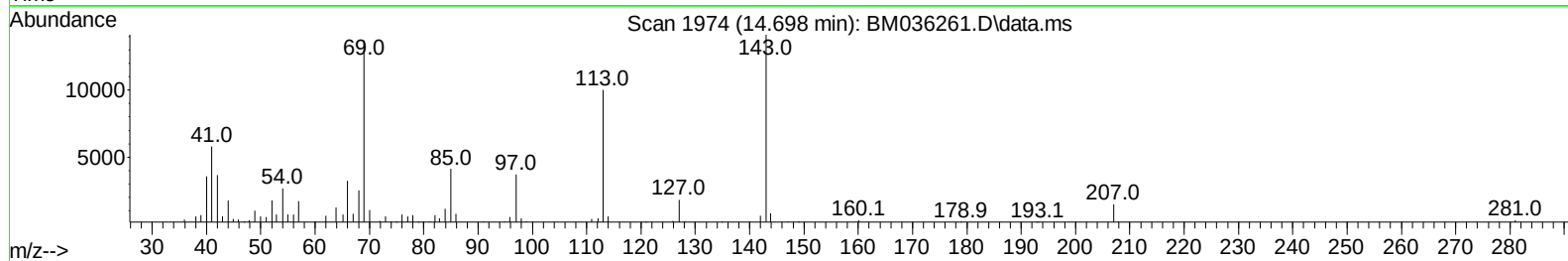
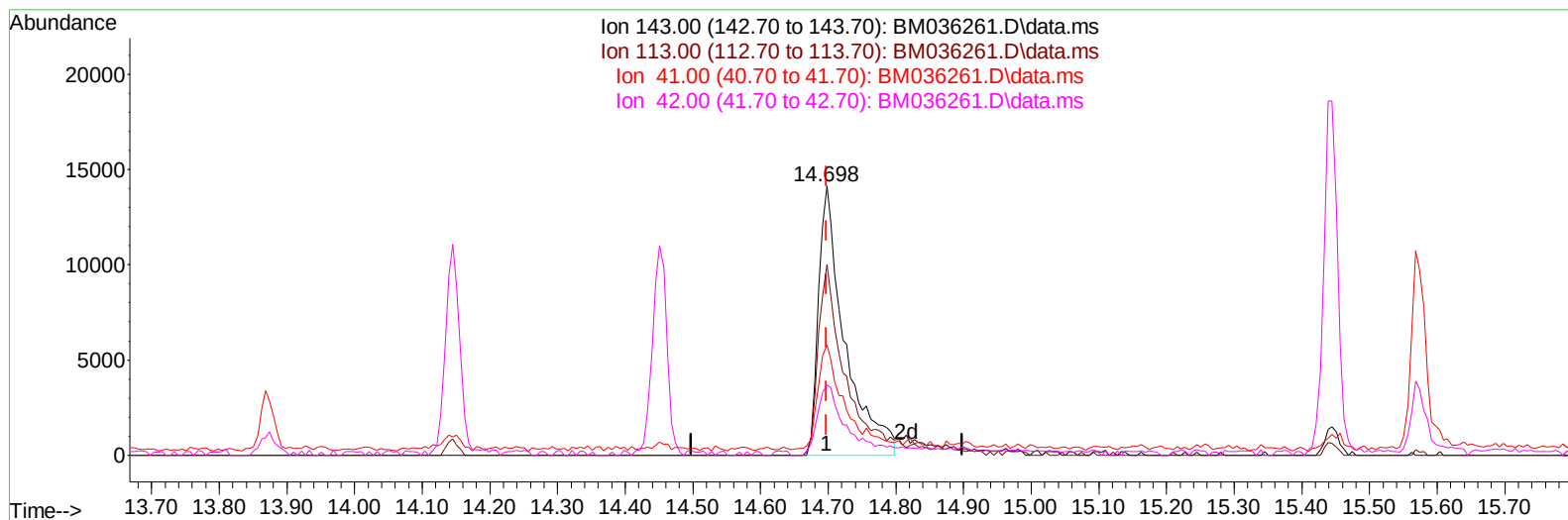
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TIC: BM036261.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.698min (-0.000) 4.62 ng/ul m

response 37768

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	70.78
41.00	41.00	40.85
42.00	29.30	26.04

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	229789	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	983996	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	584654	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	1234225	20.000	ng/ul	-0.01
79) Chrysene-d12	21.380	240	1267367	20.000	ng/ul	0.00
88) Perylene-d12	23.709	264	1348754	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	23445	3.839	ng/uL	0.00
4) Pyridine-d5	3.646	84	156655	9.364	ng/ul	0.00
7) Phenol-d5	6.998	99	118736	6.162	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.145	67	358537	27.910	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	353404	22.095	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	215633	13.813	ng/ul	-0.01
21) Nitrobenzene-d5	8.981	128	229879	29.482	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	223620	29.437	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	354652	25.812	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	565390	24.308	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	1224972	30.759	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1457645	28.686	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.698	143	37768m	4.618	ng/ul	0.00
60) Fluorene-d10	15.445	176	1152886	32.035	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	177704	26.786	ng/ul	0.00
73) Anthracene-d10	17.292	188	2007265	35.972	ng/ul	-0.01
81) Pyrene-d10	19.586	212	2369895	33.740	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	2487315	36.325	ng/ul	-0.01

Target Compounds Qvalue

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

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