

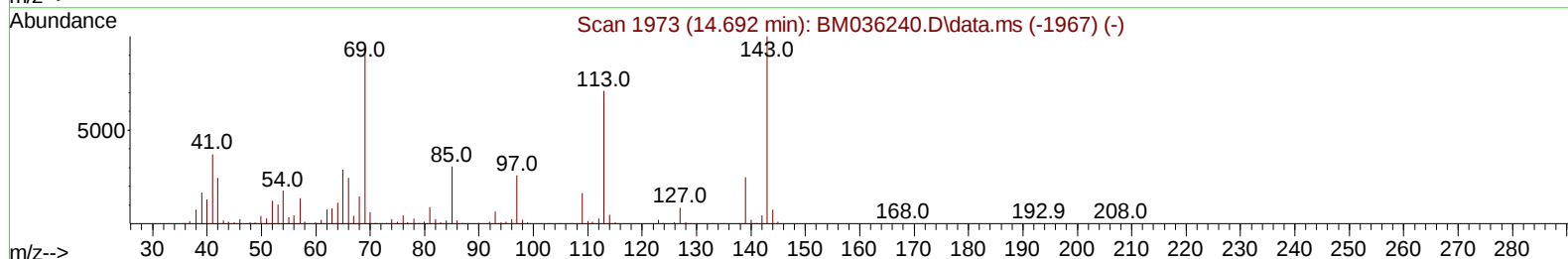
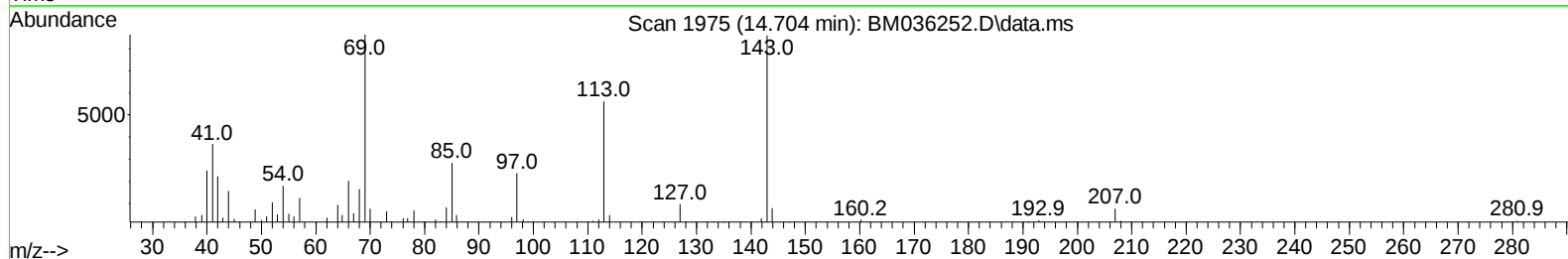
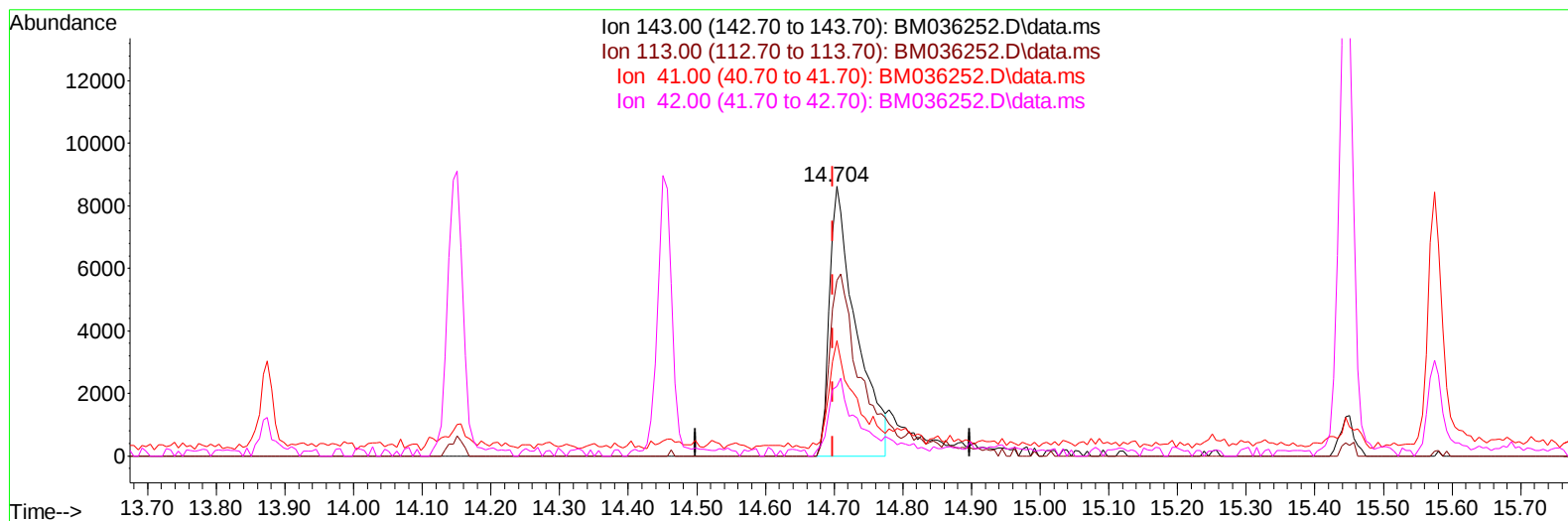
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036252.D
Acq On : 13 Aug 2022 04:58
Operator : CG/JU
Sample : N4165-05
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGBY0

Manual IntegrationsAPPROVED

Quant Time: Aug 13 06:00:41 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: BM036252.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.006) 3.40 ng/u1

response 23125

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	65.29
41.00	41.00	42.79
42.00	29.30	25.76

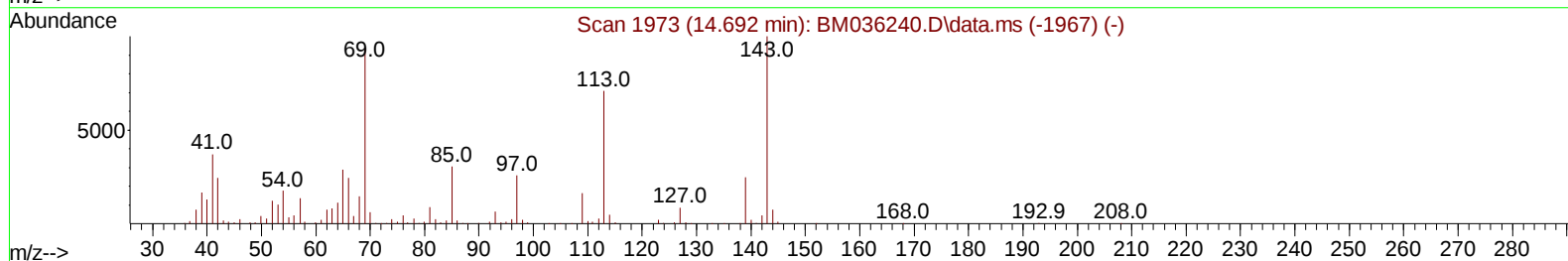
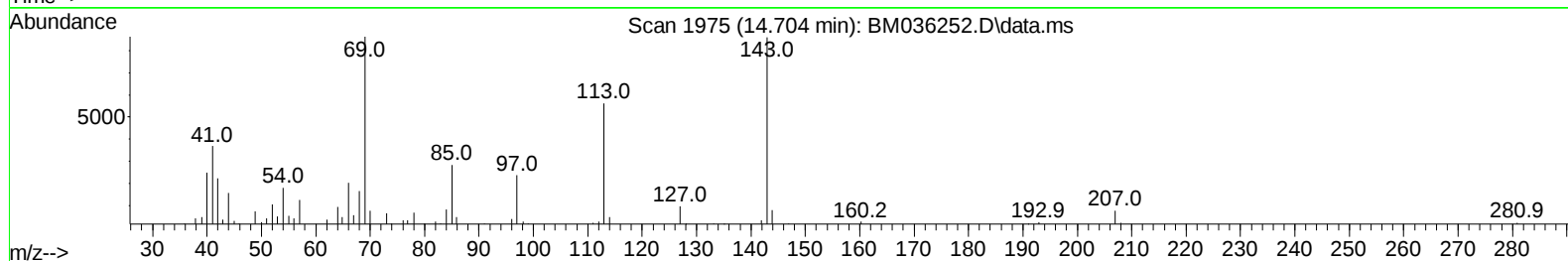
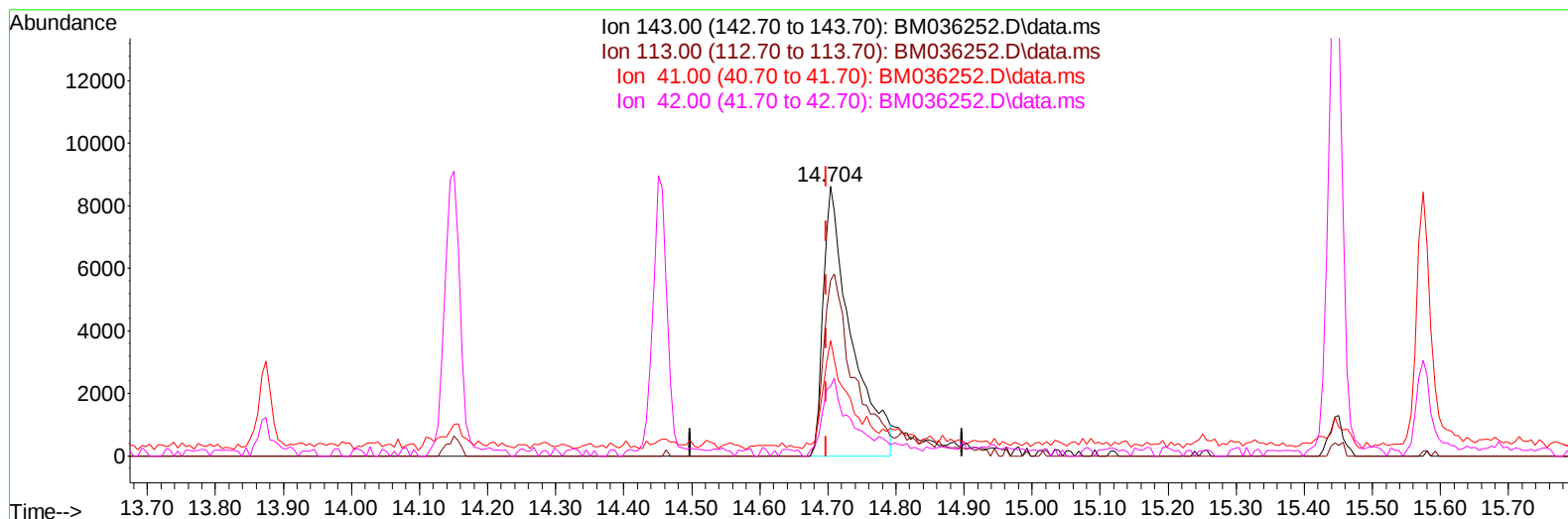
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TIC: BM036252.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.006) 3.60 ng/ul m

response 24447

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	65.29
41.00	41.00	42.79
42.00	29.30	25.76

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	209257	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	872499	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	485636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	971595	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	1002095	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1104512	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	27958	5.028	ng/uL	0.00
4) Pyridine-d5	3.646	84	144016	9.453	ng/ul	0.00
7) Phenol-d5	6.998	99	104941	5.980	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.151	67	350946	30.000	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	341410	23.439	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	191728	13.487	ng/ul	-0.01
21) Nitrobenzene-d5	8.981	128	219650	31.770	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	209471	31.098	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	328909	26.998	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	516110	25.025	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	1061235	32.080	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1299623	30.791	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	24447m	3.599	ng/ul	0.00
60) Fluorene-d10	15.445	176	985718	32.975	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	134512	25.756	ng/ul	0.00
73) Anthracene-d10	17.298	188	1615163	36.770	ng/ul	0.00
81) Pyrene-d10	19.592	212	1872995	33.724	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	2042894	36.432	ng/ul	-0.01

Target Compounds Qvalue

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

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