

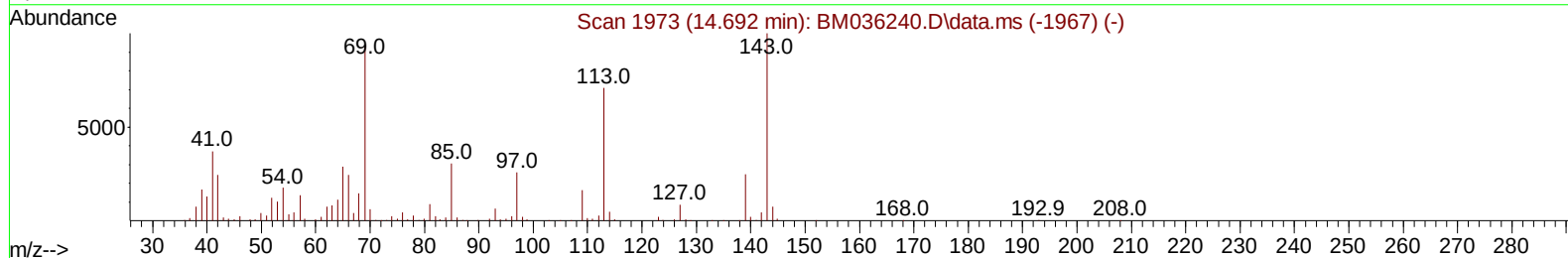
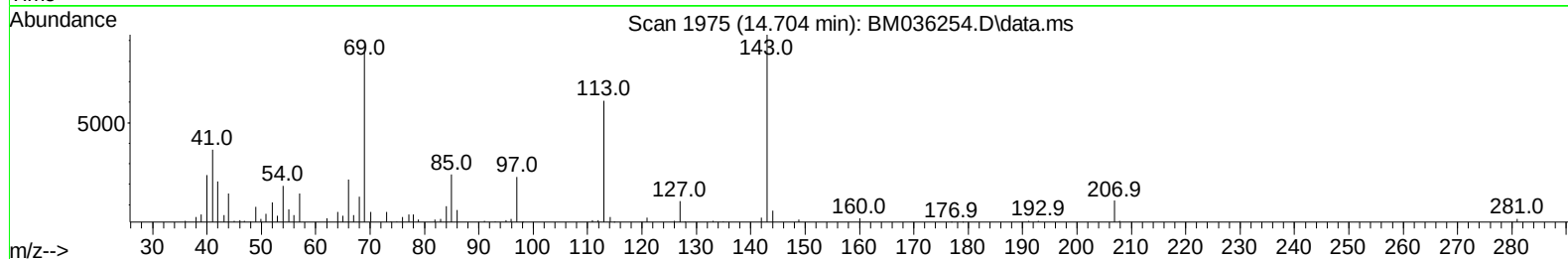
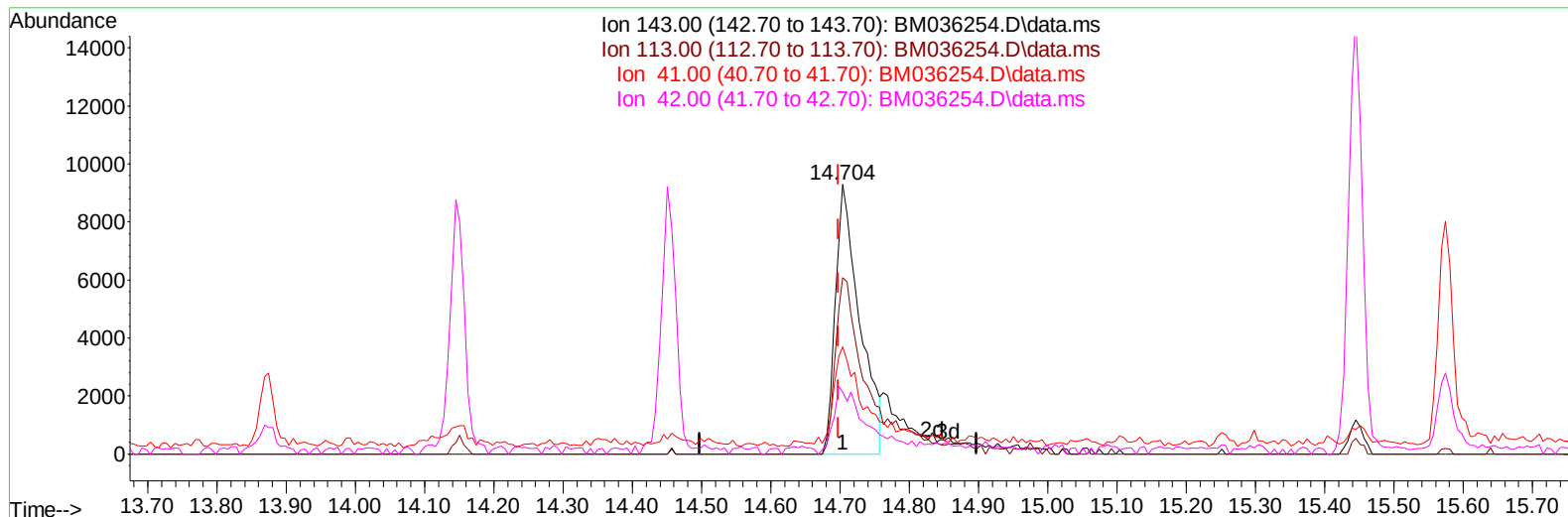
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
Data File : BM036254.D
Acq On : 13 Aug 2022 06:09
Operator : CG/JU
Sample : N4165-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGBY3

Manual IntegrationsAPPROVED

Quant Time: Aug 13 06:44:05 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: BM036254.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.006) 3.39 ng/ul

response 22551

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	65.49
41.00	41.00	39.81
42.00	29.30	22.90#

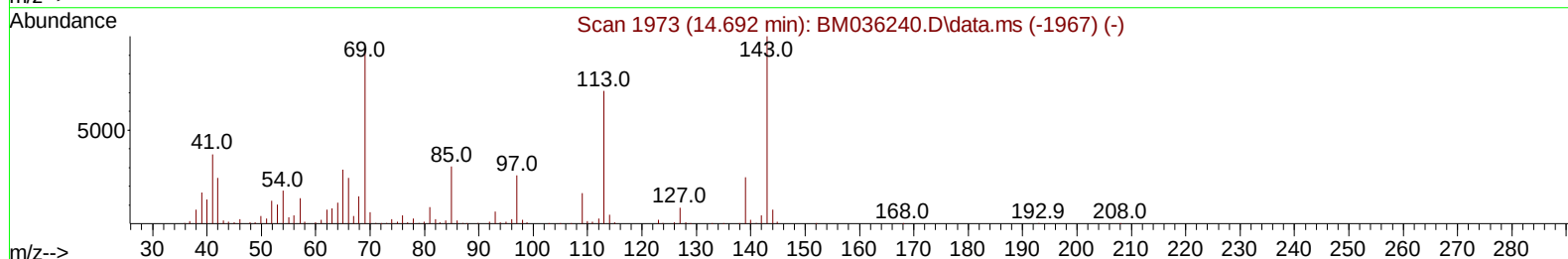
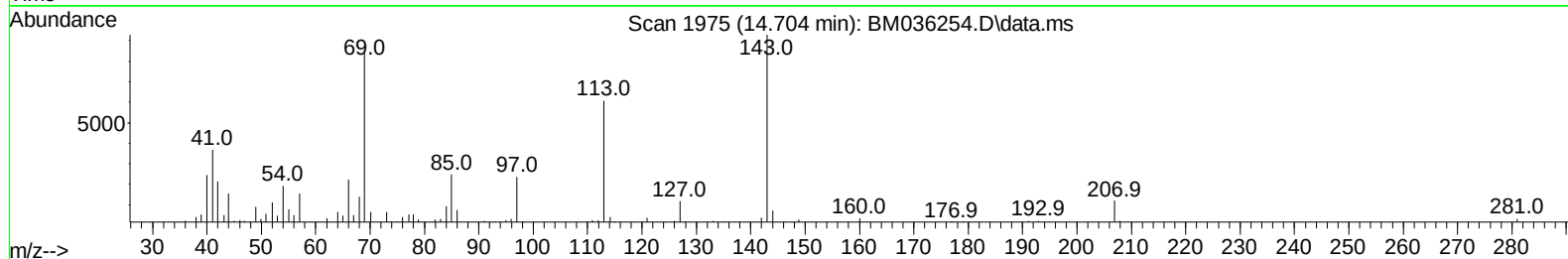
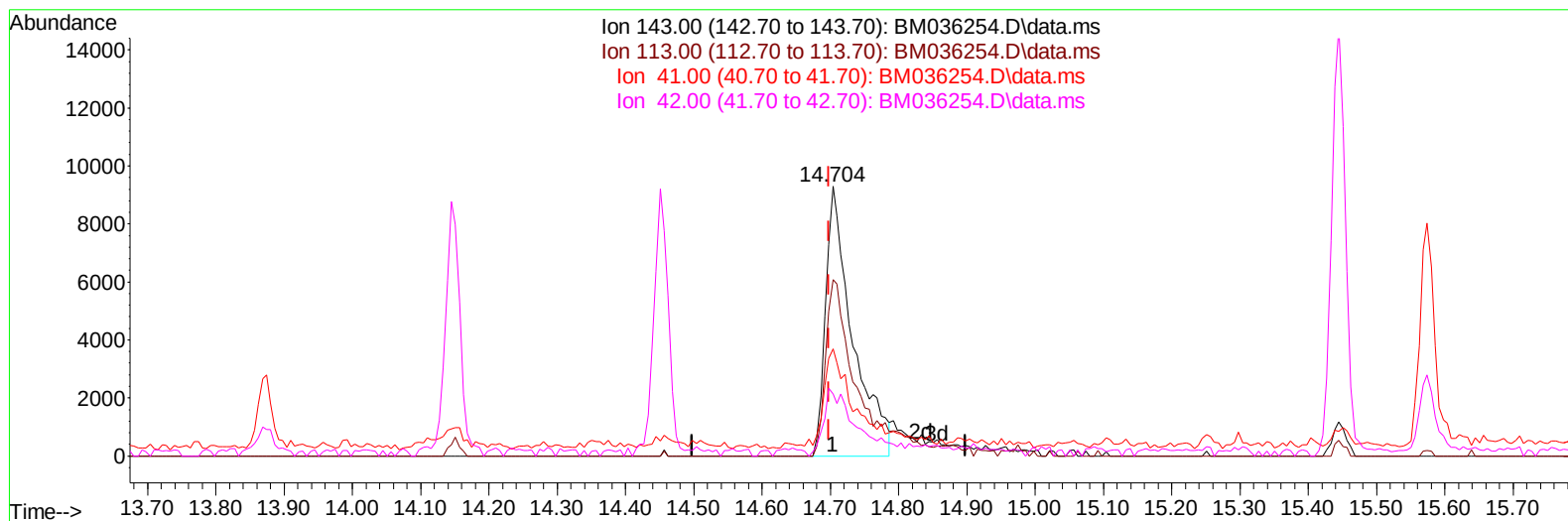
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TIC: BM036254.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.006) 3.82 ng/ul m

response 25365

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	65.49
41.00	41.00	39.81
42.00	29.30	22.90#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	200977	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	834485	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	475156	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	996636	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	1096385	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1247050	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	25282	4.734	ng/uL	0.00
4) Pyridine-d5	3.652	84	96097	6.568	ng/ul	0.00
7) Phenol-d5	6.998	99	91344	5.420	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.151	67	306603	27.289	ng/ul	0.00
11) 2-Chlorophenol-d4	7.345	132	298470	21.335	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	166521	12.196	ng/ul	-0.01
21) Nitrobenzene-d5	8.981	128	188747	28.544	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	181898	28.234	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	288430	24.754	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	377049	19.115	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	981674	30.330	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1186962	28.742	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.704	143	25365m	3.816	ng/ul	0.00
60) Fluorene-d10	15.445	176	918558	31.406	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	132400	24.714	ng/ul	0.00
73) Anthracene-d10	17.292	188	1591538	35.322	ng/ul	-0.01
81) Pyrene-d10	19.592	212	1907570	31.393	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	2223108	35.114	ng/ul	-0.01

Target Compounds Qvalue

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

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