

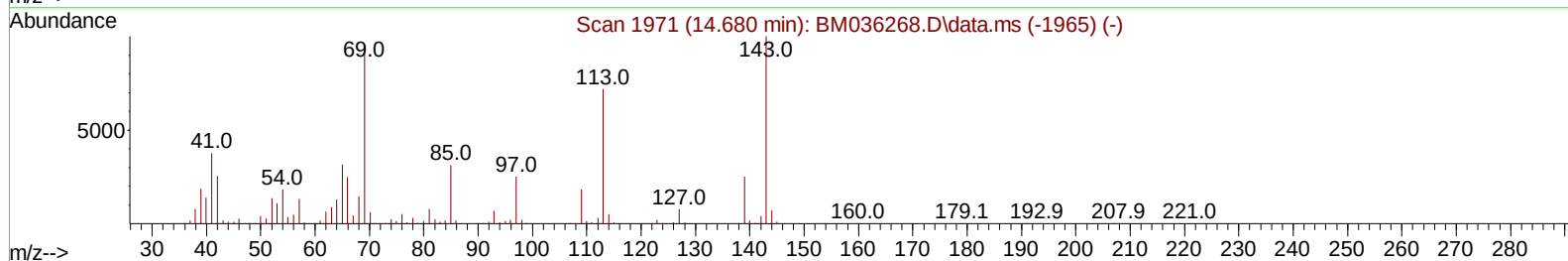
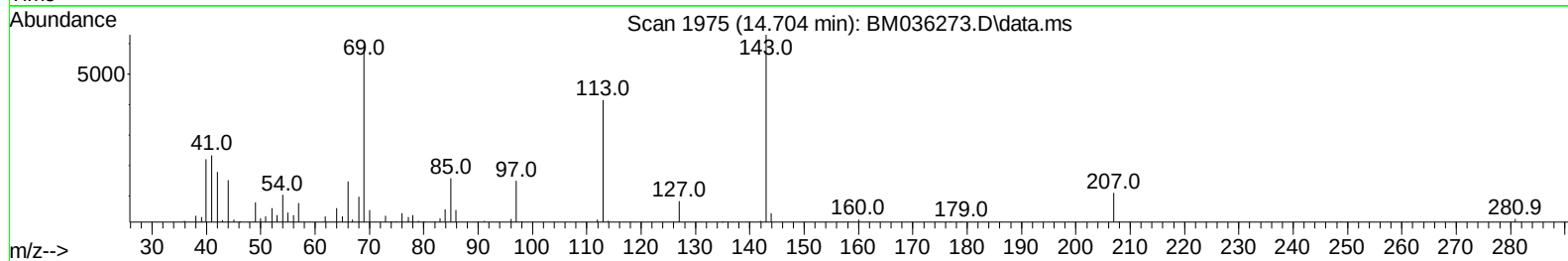
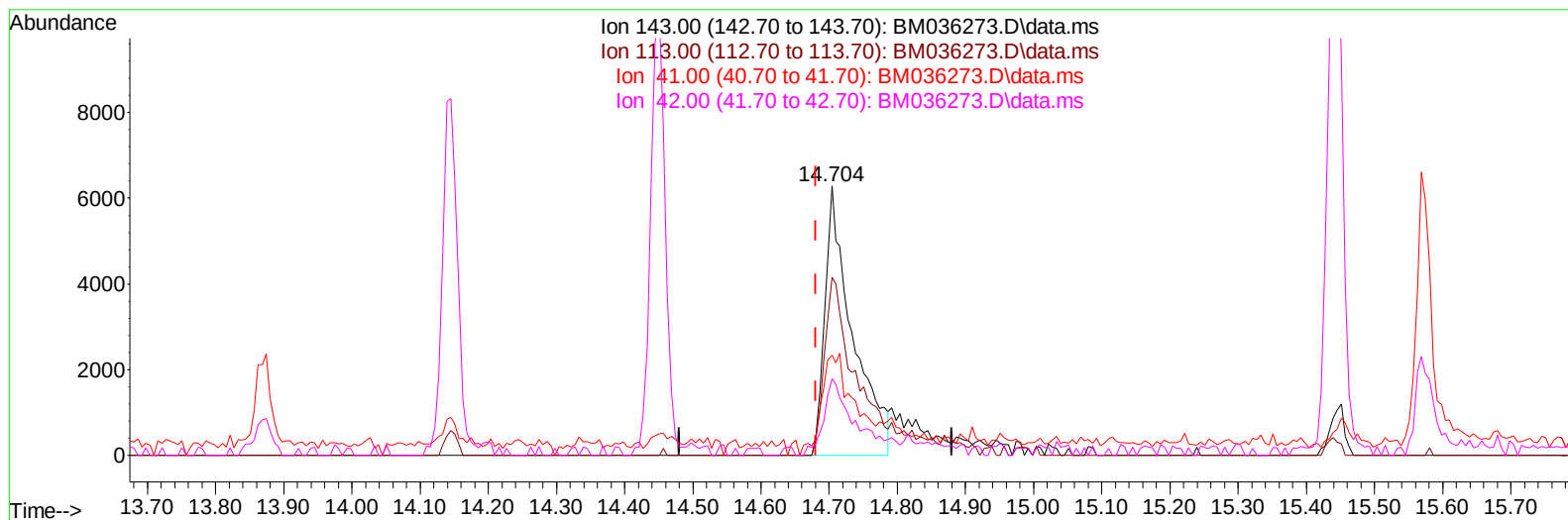
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081422\
 Data File : BM036273.D
 Acq On : 14 Aug 2022 16:43
 Operator : CG/JU
 Sample : N4166-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGC16

Manual IntegrationsAPPROVED

Quant Time: Aug 14 23:38:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Aug 14 23:34:05 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2022
 Supervised By :mohammad ahmed 08/16/2022



TIC: BM036273.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.023) 2.37 ng/ul

response 17602

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	66.10
41.00	41.00	37.20
42.00	29.30	28.44

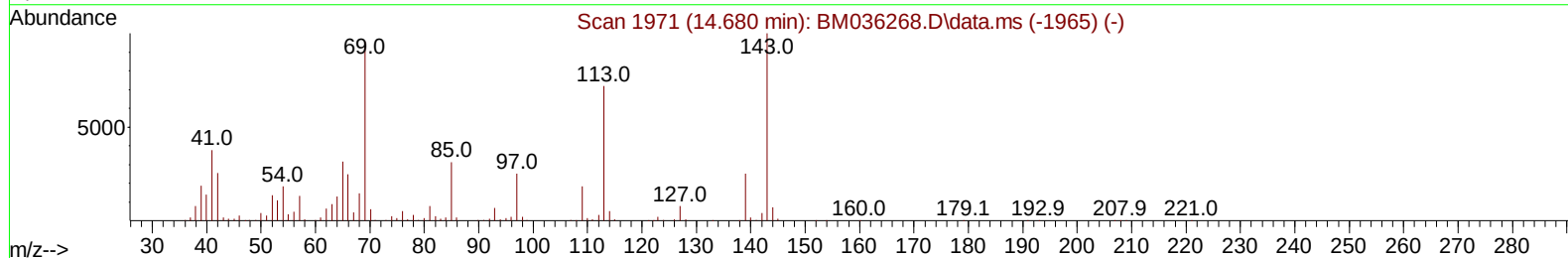
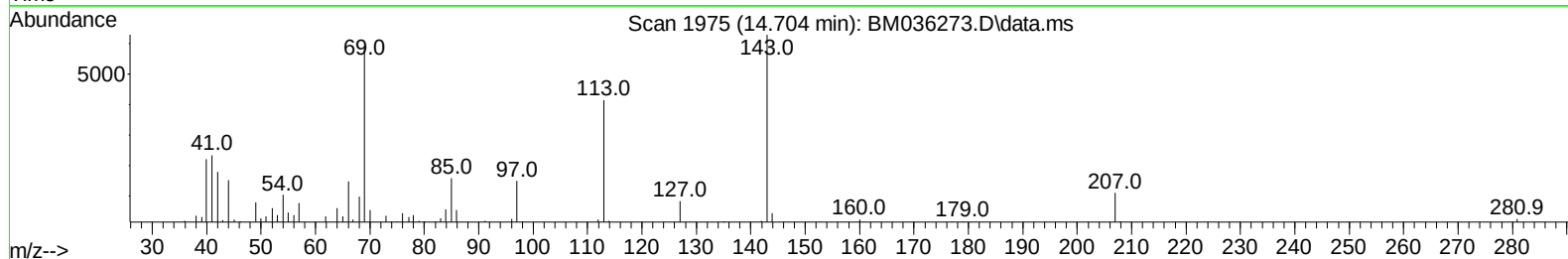
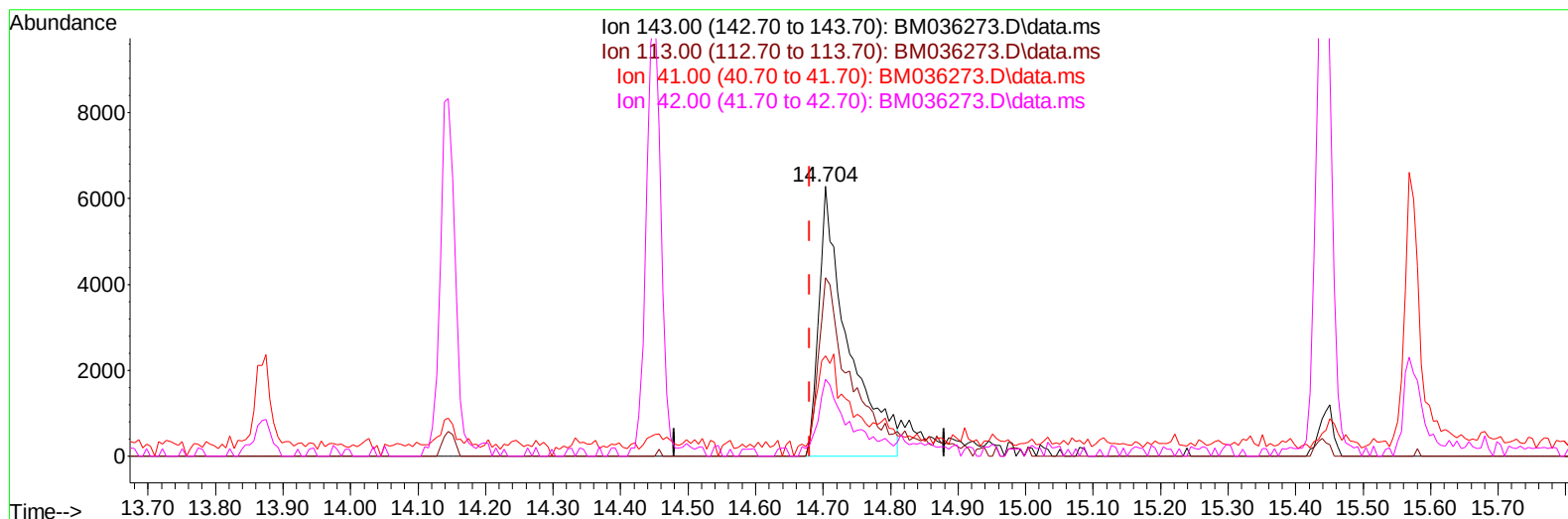
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(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.023) 2.53 ng/ul m

response 18842

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	66.10
41.00	41.00	37.20
42.00	29.30	28.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	228831	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	949829	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	531818	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	1023638	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	950504	20.000	ng/ul	0.00
88) Perylene-d12	23.709	264	1000722	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	23428	3.853	ng/uL	0.00
4) Pyridine-d5	3.646	84	102809	6.171	ng/ul	0.00
7) Phenol-d5	6.992	99	88538	4.614	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	300101	23.459	ng/ul	0.00
11) 2-Chlorophenol-d4	7.340	132	286964	18.016	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	164636	10.590	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	182935	24.305	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	175159	23.887	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.239	165	274548	20.701	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	420335	18.722	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	939080	25.923	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1101230	23.825	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	18842m	2.533	ng/ul	0.02
60) Fluorene-d10	15.445	176	861573	26.319	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	110805	20.138	ng/ul	0.00
73) Anthracene-d10	17.292	188	1411910	30.508	ng/ul	0.00
81) Pyrene-d10	19.586	212	1550852	29.440	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.556	264	1551612	30.540	ng/ul	0.00

Target Compounds Qvalue

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

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