

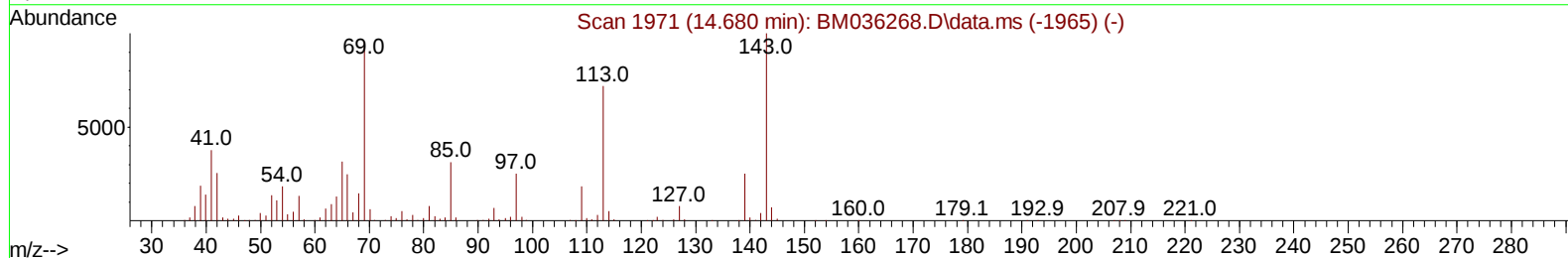
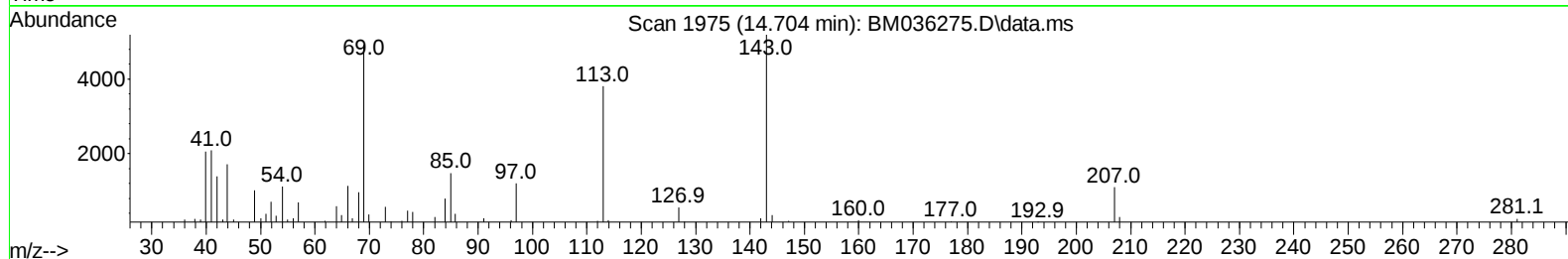
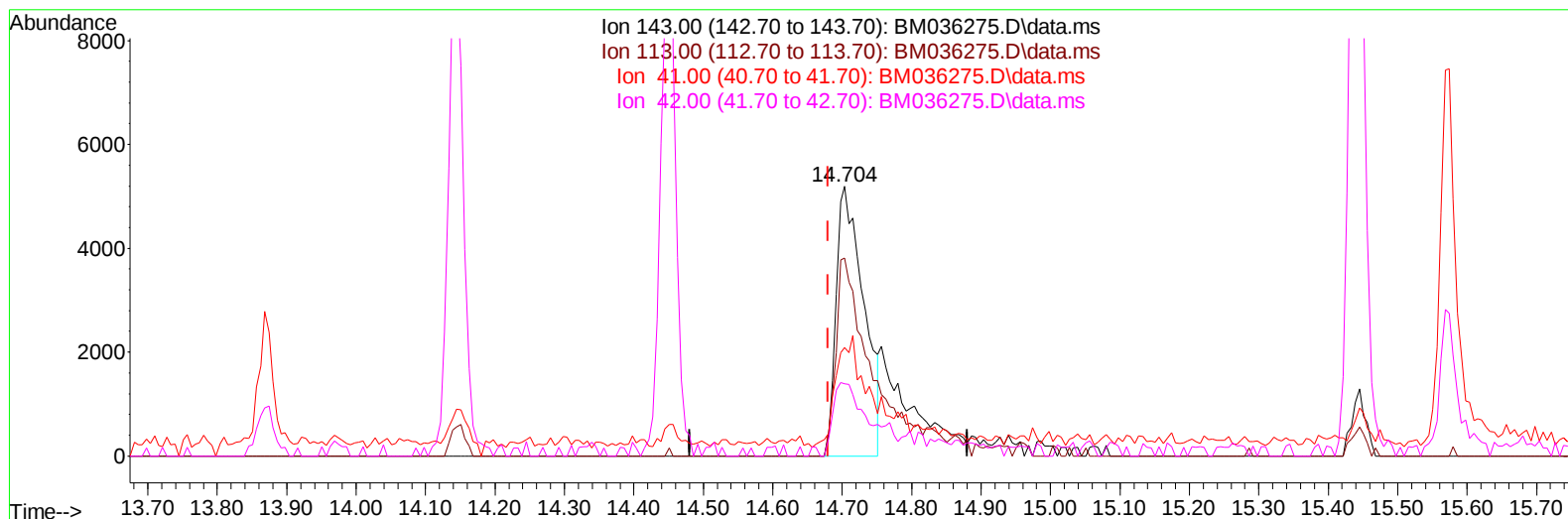
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081422\
Data File : BM036275.D
Acq On : 14 Aug 2022 17:56
Operator : CG/JU
Sample : N4166-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BGC02

Manual IntegrationsAPPROVED

Quant Time: Aug 14 23:39:14 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: BM036275.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.023) 2.04 ng/ul

response 14220

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	73.55
41.00	41.00	40.17
42.00	29.30	26.67

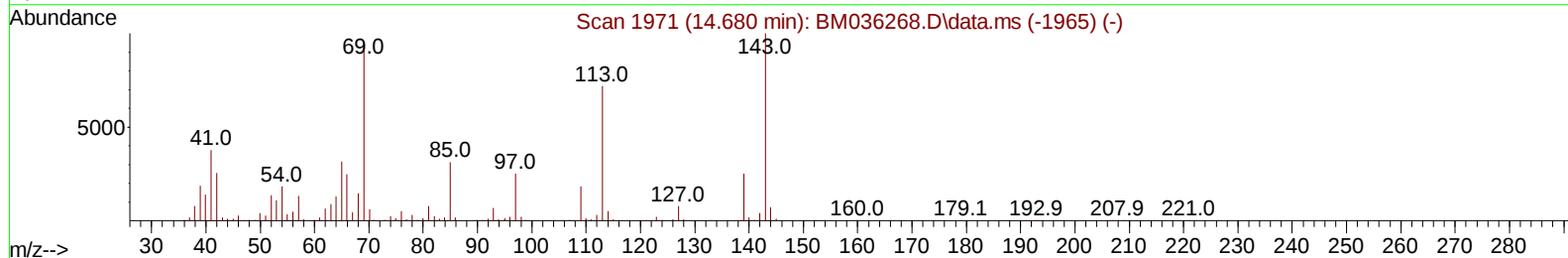
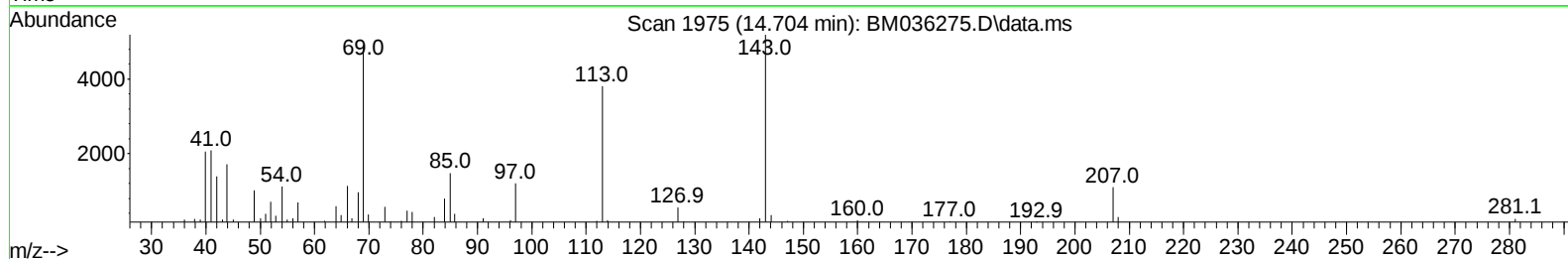
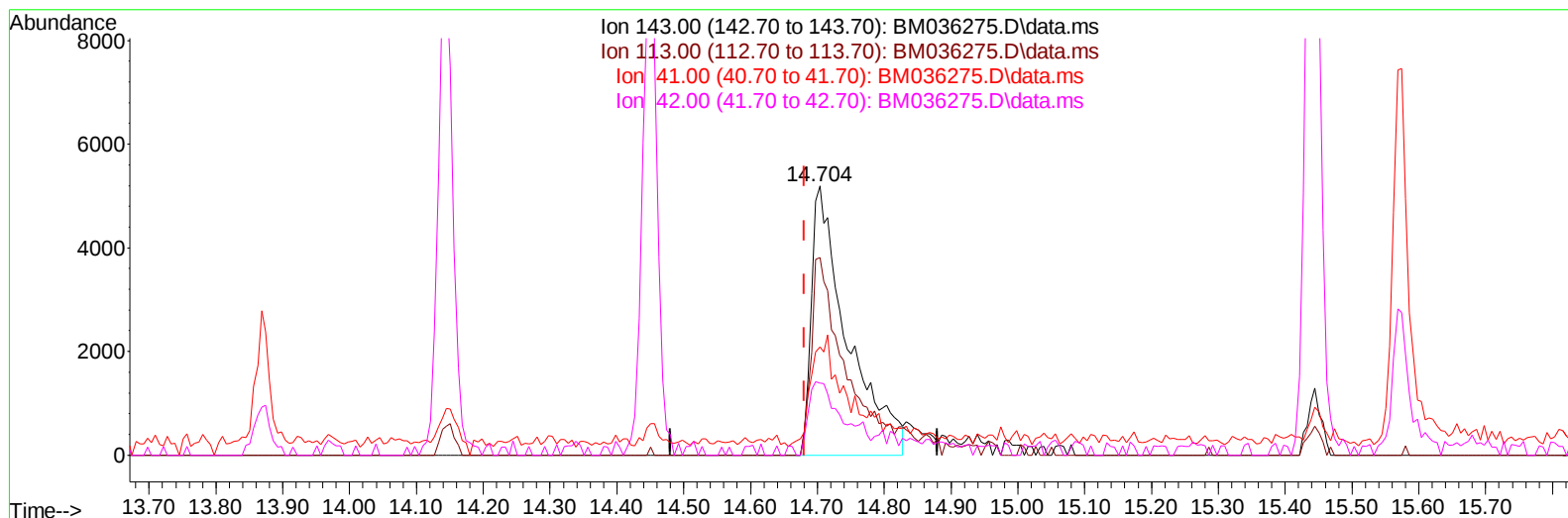
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TIC: BM036275.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (+ 0.023) 2.77 ng/ul m

response 19307

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	73.55
41.00	41.00	40.17
42.00	29.30	26.67

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	212059	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	873916	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	498287	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	990219	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	984801	20.000	ng/ul	0.00
88) Perylene-d12	23.709	264	1079638	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	25229	4.477	ng/uL	0.00
4) Pyridine-d5	3.651	84	65052	4.214	ng/ul	0.01
7) Phenol-d5	6.998	99	81919	4.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	319803	26.976	ng/ul	0.00
11) 2-Chlorophenol-d4	7.339	132	289179	19.591	ng/ul	0.00
15) 4-Methylphenol-d8	8.533	113	157863	10.958	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	197340	28.497	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	183241	27.160	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.239	165	290601	23.815	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	413364	20.010	ng/ul	0.00
46) Dimethylphthalate-d6	13.868	166	1020632	30.070	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1232644	28.463	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	19307m	2.770	ng/ul	0.02
60) Fluorene-d10	15.445	176	947669	30.897	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	126214	23.712	ng/ul	0.00
73) Anthracene-d10	17.292	188	1601754	35.779	ng/ul	0.00
81) Pyrene-d10	19.586	212	1789889	32.794	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.556	264	1907089	34.793	ng/ul	0.00

Target Compounds Qvalue

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

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