

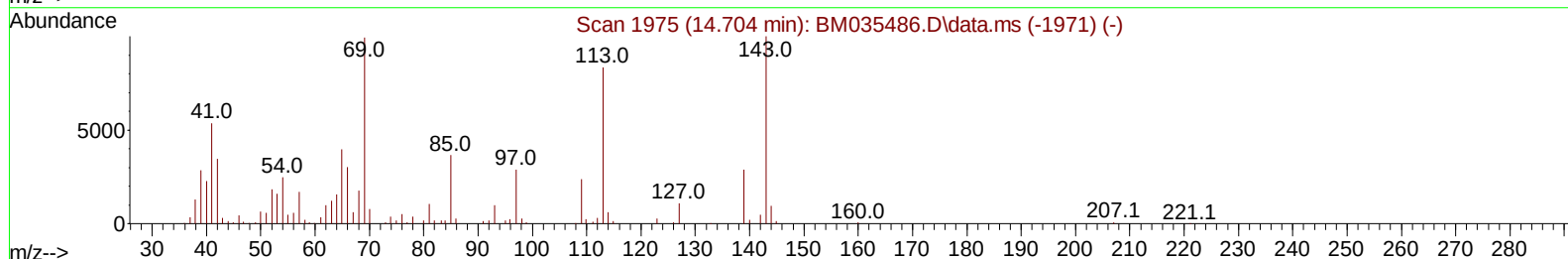
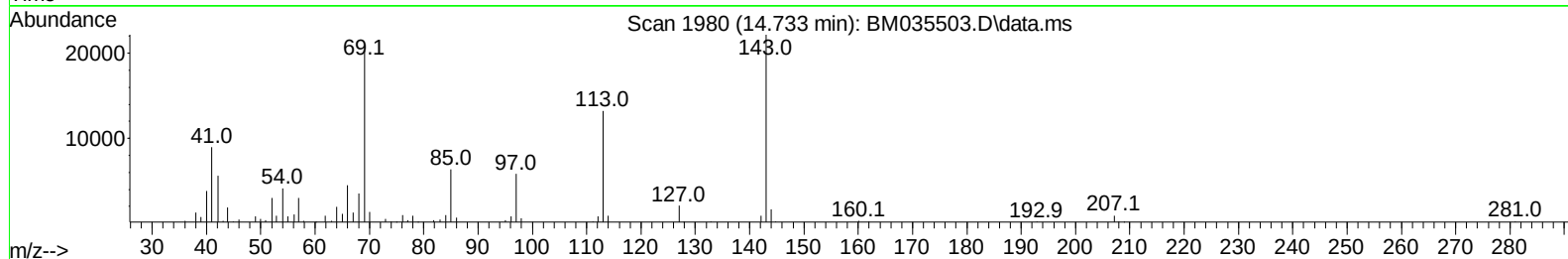
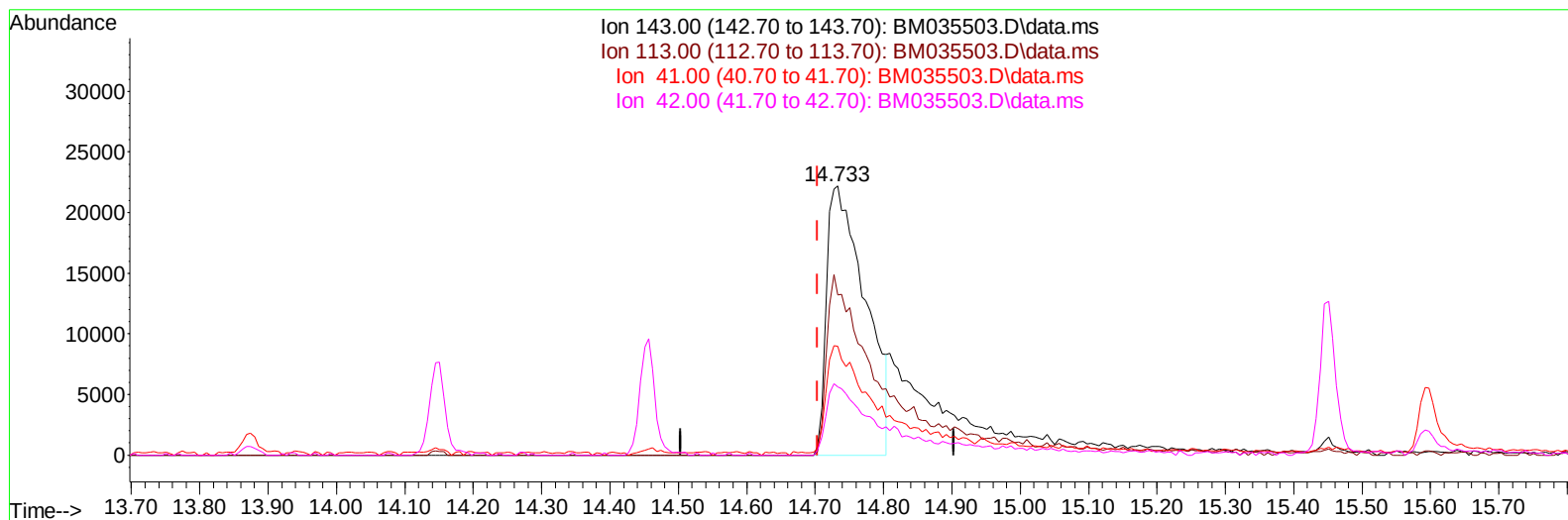
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
Data File : BM035503.D
Acq On : 17 Jun 2022 20:45
Operator : CG/JU
Sample : PB145488BL
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK488

Manual IntegrationsAPPROVED

Quant Time: Jun 17 23:03:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 01:58:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/18/2022
Supervised By :Yogesh Patel 06/22/2022



TIC: BM035503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.029) 16.00 ng/ul

response 87184

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	59.56
41.00	41.00	40.38
42.00	29.30	25.46

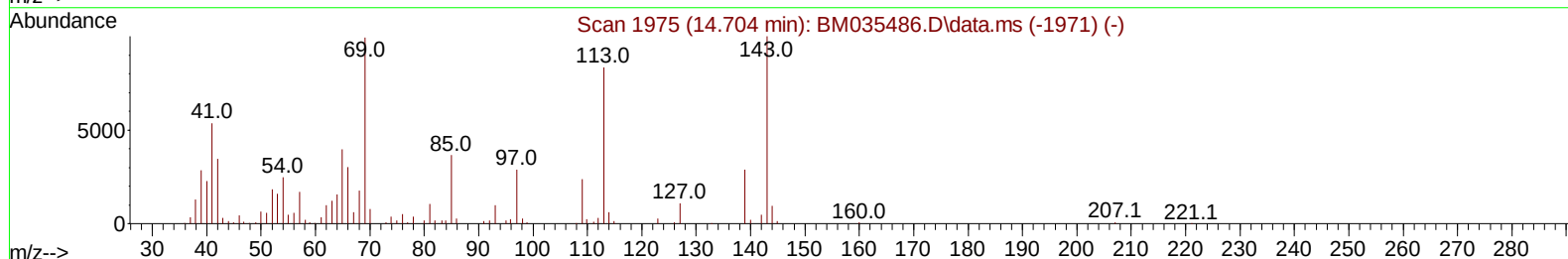
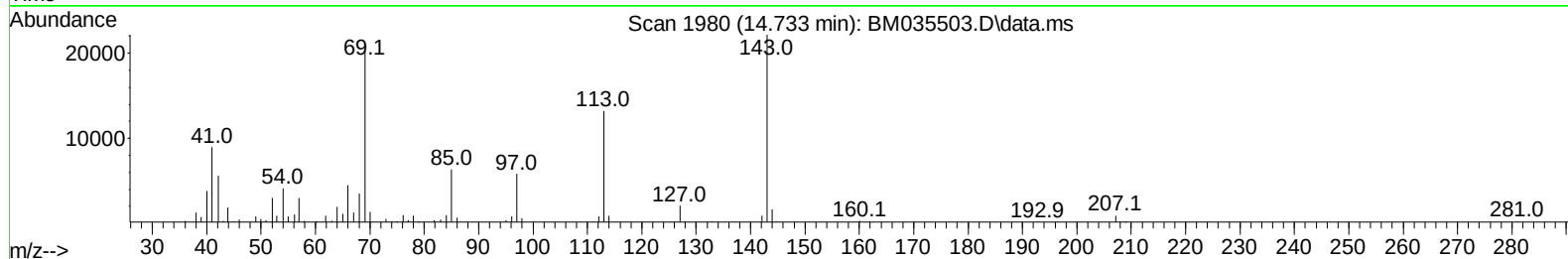
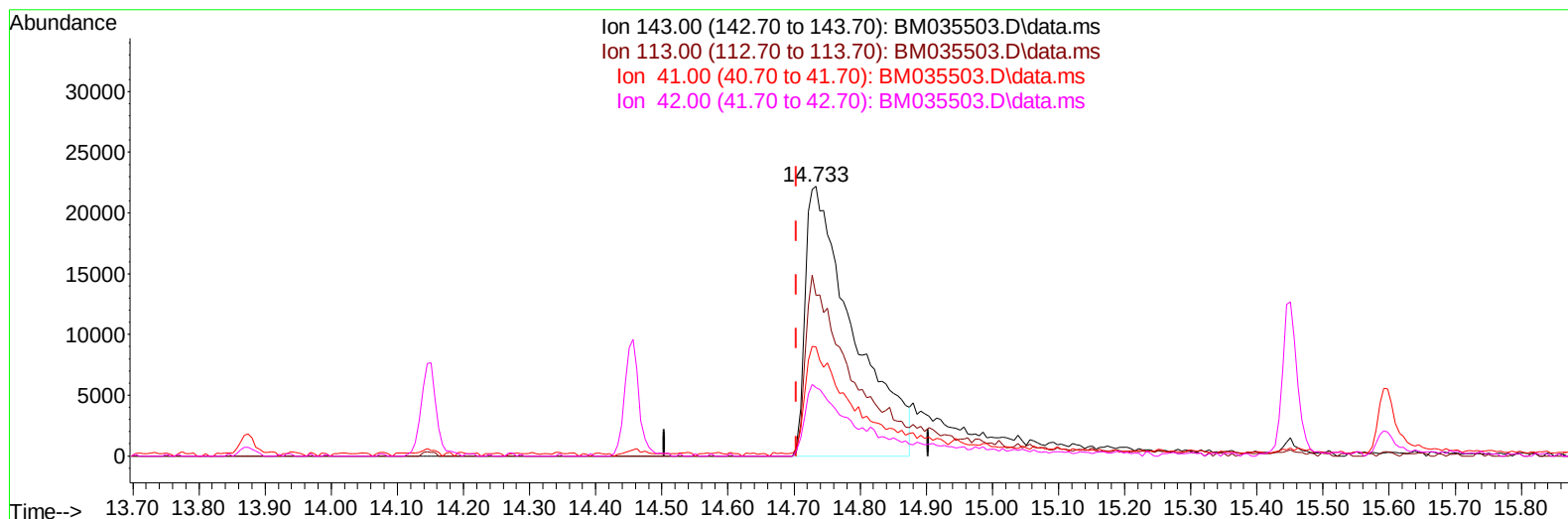
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TIC: BM035503.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.029) 20.50 ng/ul m

response 111722

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	59.56
41.00	41.00	40.38
42.00	29.30	25.46

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	209666	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	886917	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	521642	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.203	188	1061220	20.000	ng/ul	0.00
79) Chrysene-d12	21.391	240	1070547	20.000	ng/ul	0.00
88) Perylene-d12	23.727	264	1192304	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	26955	5.132	ng/uL	0.00
4) Pyridine-d5	3.687	84	275128	21.268	ng/ul	0.00
7) Phenol-d5	7.004	99	387021	24.085	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	247707	24.789	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	368394	26.469	ng/ul	0.00
15) 4-Methylphenol-d8	8.539	113	331890	24.673	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	162520	24.617	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	180574	25.365	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	268656	22.227	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	692359	36.405	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	974880	27.475	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1156920	26.311	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	111722m	20.497	ng/ul	0.03
60) Fluorene-d10	15.451	176	850509	27.683	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	122727	19.398	ng/ul	0.00
73) Anthracene-d10	17.304	188	1362476	28.563	ng/ul	0.00
81) Pyrene-d10	19.597	212	1555912	29.562	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264	1698948	29.543	ng/ul	0.00

Target Compounds Qvalue

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

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