

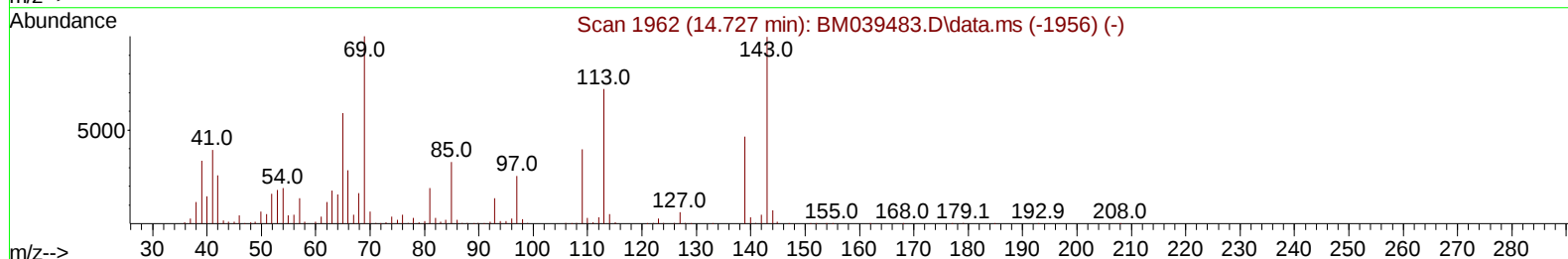
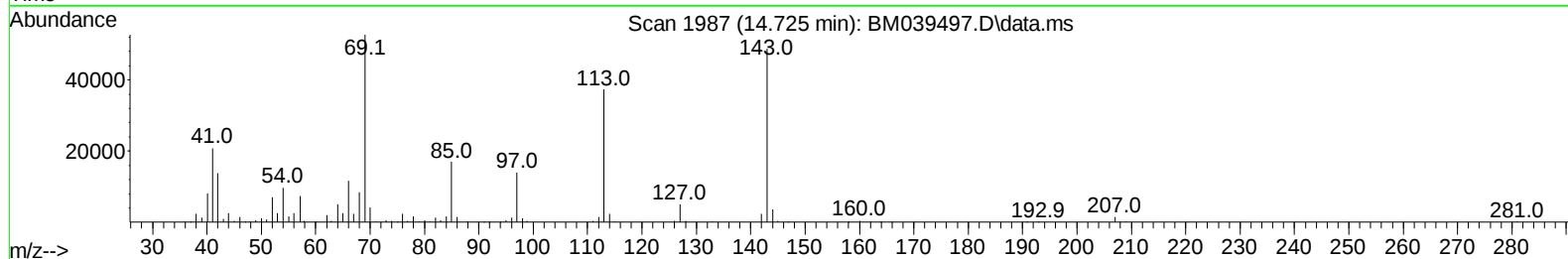
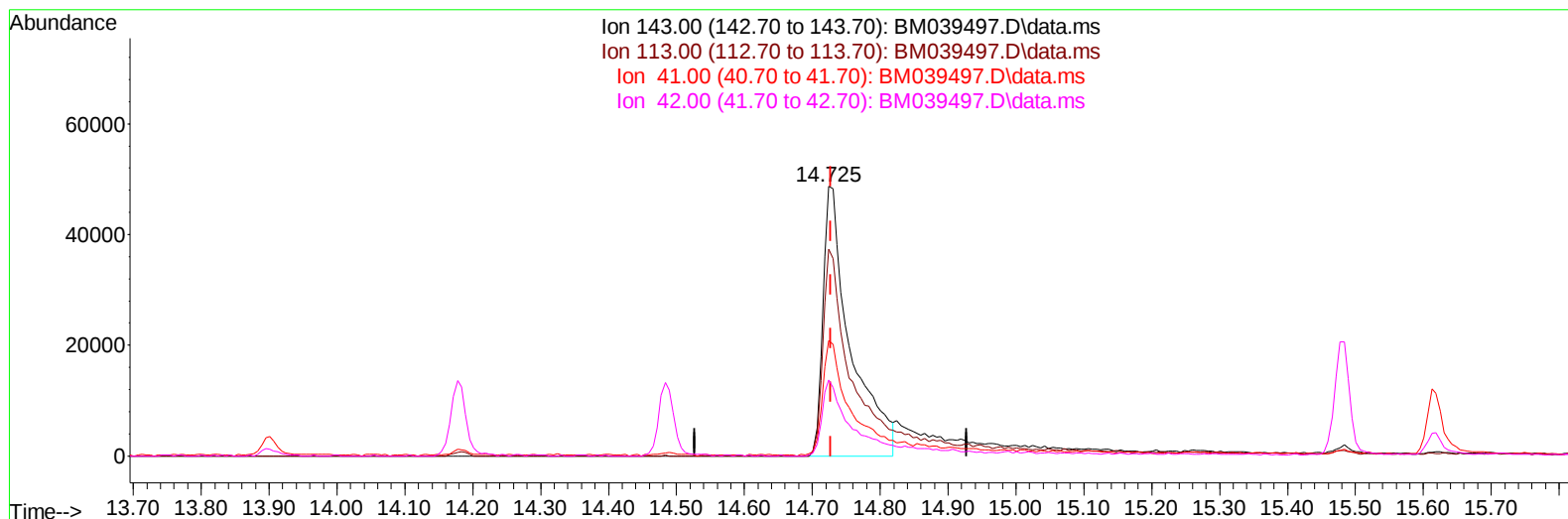
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041823\
Data File : BM039497.D
Acq On : 19 Apr 2023 16:16
Operator : CG/JU
Sample : PB152215BL
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK215

Manual IntegrationsAPPROVED

Quant Time: Apr 19 23:39:51 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 19 08:18:41 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/20/2023
Supervised By :Jagrut Upadhyay 04/20/2023



TIC: BM039497.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.725min (-0.003) 17.97 ng/u1

response 136320

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	76.63
41.00	39.90	42.80
42.00	26.60	28.29

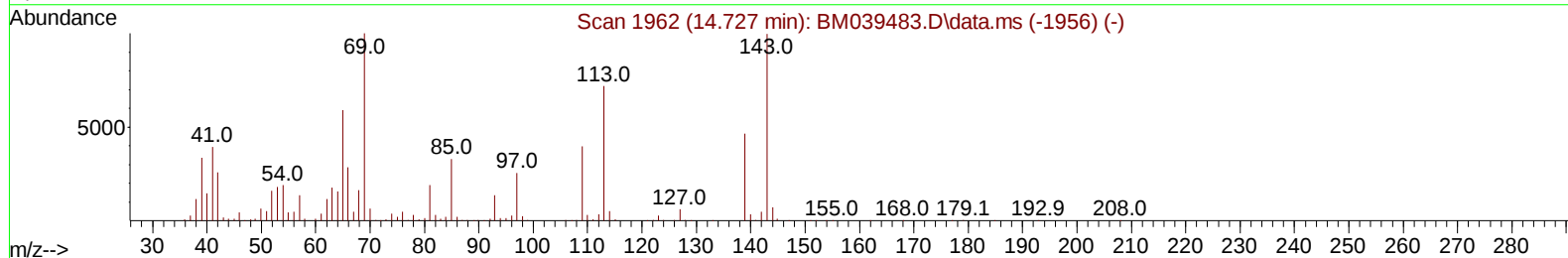
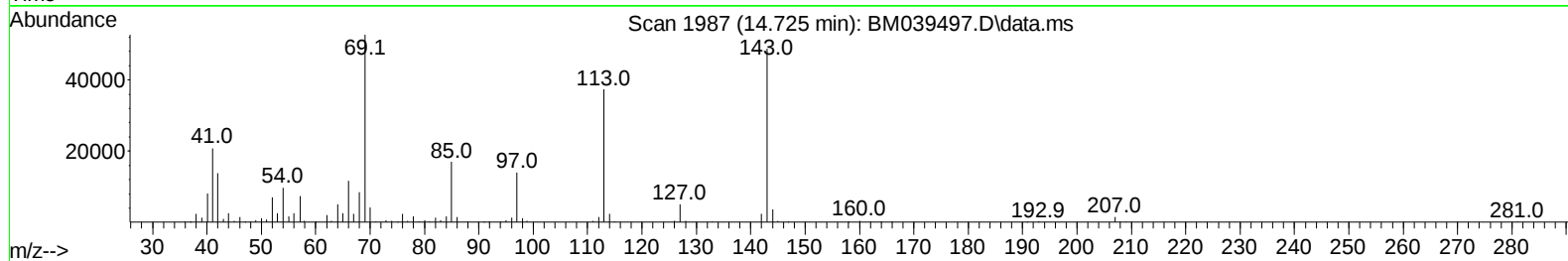
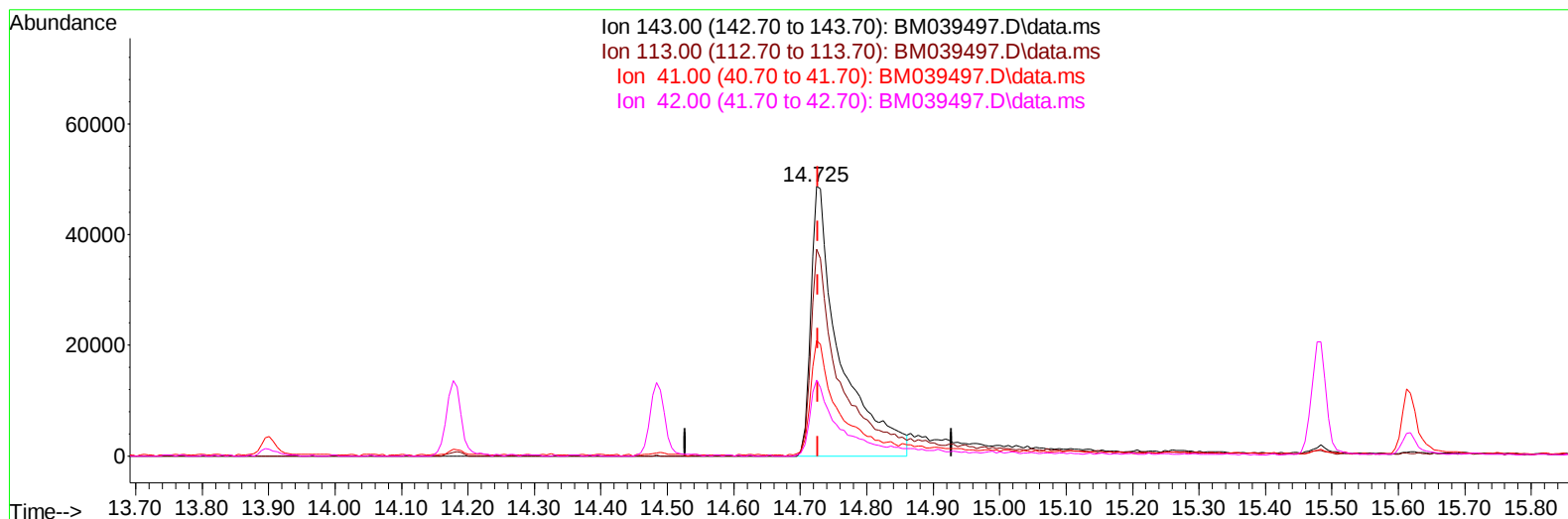
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TIC: BM039497.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.725min (-0.003) 19.56 ng/ul m

response 148356

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	76.63
41.00	39.90	42.80
42.00	26.60	28.29

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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/20/2023
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Quant Time: Apr 19 23:40:14 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.831	152	197301	20.000	ng/ul	0.00
20) Naphthalene-d8	10.637	136	918102	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.483	164	621406	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.236	188	1339822	20.000	ng/ul	0.00
79) Chrysene-d12	21.430	240	1157897	20.000	ng/ul	0.00
88) Perylene-d12	23.800	264	1087928	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	31307	6.091	ng/uL	0.00
4) Pyridine-d5	3.643	84	385359	26.682	ng/ul	0.00
7) Phenol-d5	7.013	99	474449	26.274	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.160	67	352515	29.535	ng/ul	0.00
11) 2-Chlorophenol-d4	7.360	132	405968	28.659	ng/ul	0.00
15) 4-Methylphenol-d8	8.560	113	392567	26.988	ng/ul	0.00
21) Nitrobenzene-d5	9.001	128	207436	28.084	ng/ul	0.00
24) 2-Nitrophenol-d4	9.725	143	231213	27.877	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.272	165	362483	25.400	ng/ul	0.00
31) 4-Chloroaniline-d4	10.789	131	547963	25.708	ng/ul	0.00
46) Dimethylphthalate-d6	13.901	166	1505137	30.033	ng/ul	0.00
49) Acenaphthylene-d8	14.177	160	1630489	28.960	ng/ul	0.00
54) 4-Nitrophenol-d4	14.725	143	148356m	19.560	ng/ul	0.00
60) Fluorene-d10	15.483	176	1248150	30.399	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.619	200	207431	23.298	ng/ul	0.00
73) Anthracene-d10	17.336	188	1978880	30.711	ng/ul	0.00
81) Pyrene-d10	19.636	212	2288504	29.743	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.647	264	1854250	31.569	ng/ul	0.00

Target Compounds Qvalue

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

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