

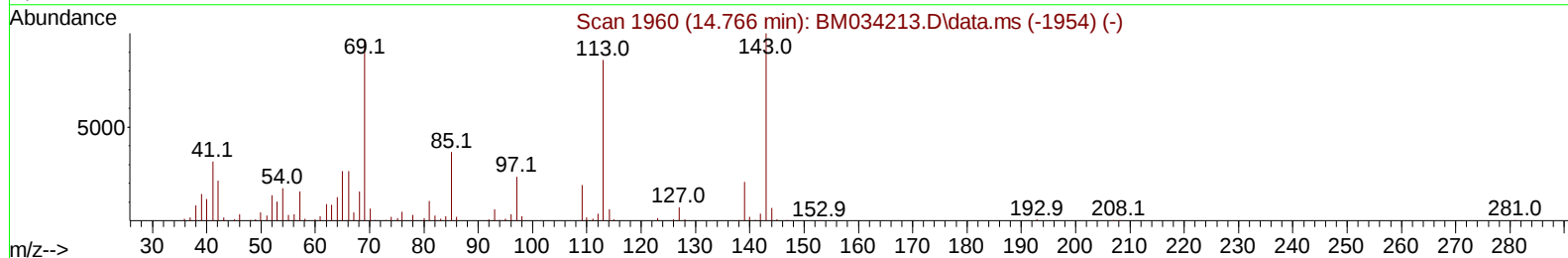
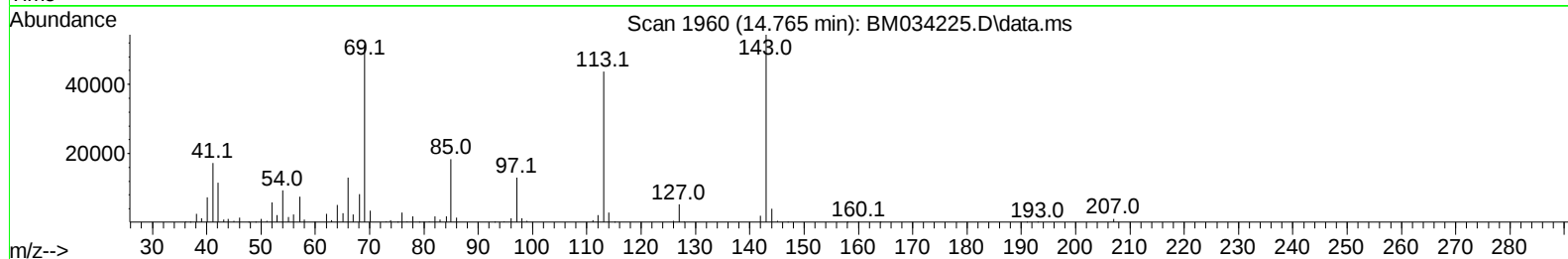
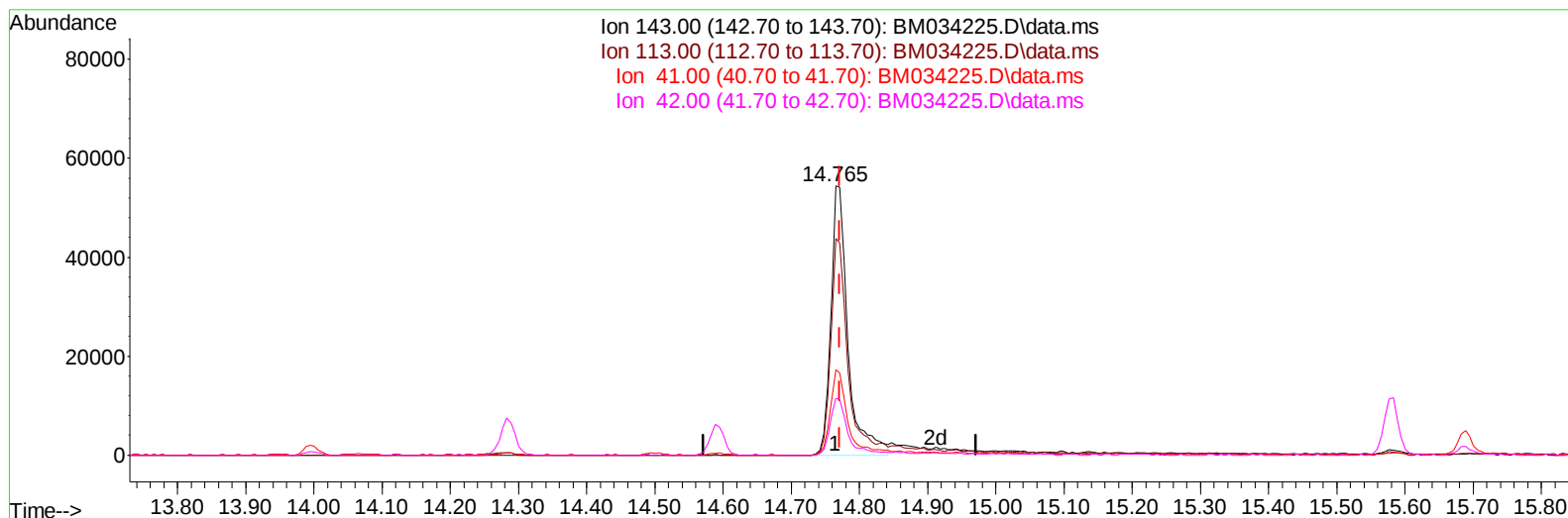
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
 Data File : BM034225.D
 Acq On : 11 Mar 2022 16:47
 Operator : CG/JU
 Sample : N1816-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSD2

Manual IntegrationsAPPROVED

Quant Time: Mar 11 23:41:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 11 02:00:05 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/14/2022
 Supervised By :mohammad ahmed 03/14/2022



TIC: BM034225.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (-0.006) 23.89 ng/u1

response 94902

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	80.43
41.00	31.90	31.81
42.00	19.10	21.15

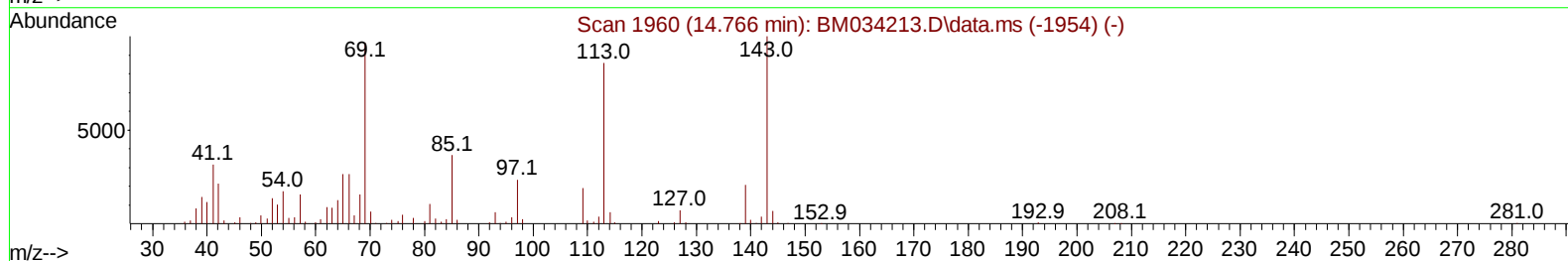
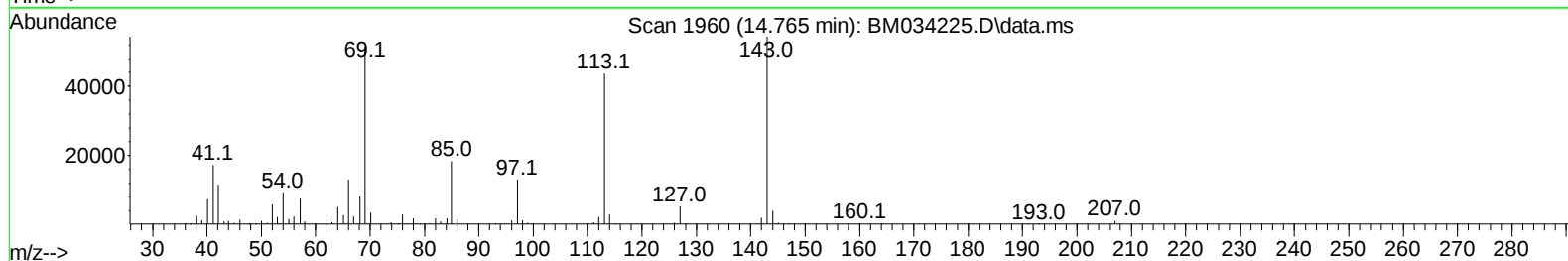
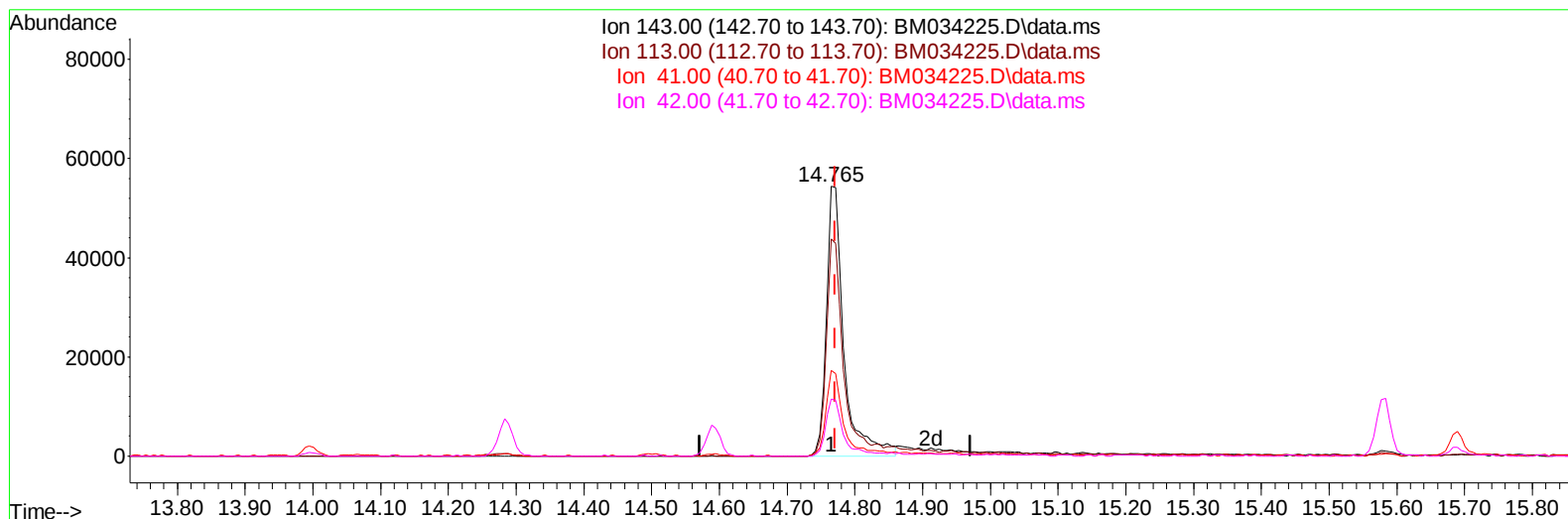
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(54) 4-Nitrophenol-d4 (S)

14.765min (-0.006) 24.47 ng/ul m

response 97211

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	80.43
41.00	31.90	31.81
42.00	19.10	21.15

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	135247	20.000	ng/ul	0.00
20) Naphthalene-d8	10.766	136	528673	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.595	164	350546	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	703437	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	651734	20.000	ng/ul	0.00
88) Perylene-d12	23.924	264	703217	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	20716	6.686	ng/uL	0.00
4) Pyridine-d5	3.743	84	250238	33.923	ng/ul	0.00
7) Phenol-d5	7.101	99	308816	34.120	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	198658	37.414	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	292160	37.351	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	268265	34.656	ng/ul	0.00
21) Nitrobenzene-d5	9.113	128	141914	37.423	ng/ul	0.00
24) 2-Nitrophenol-d4	9.842	143	153318	36.461	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.377	165	298782	35.443	ng/ul	0.00
31) 4-Chloroaniline-d4	10.895	131	362098	34.539	ng/ul	0.00
46) Dimethylphthalate-d6	14.001	166	953928	37.400	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	1191447	36.778	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	97211m	24.468	ng/ul	0.00
60) Fluorene-d10	15.583	176	839892	37.617	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	120503	27.076	ng/ul	0.00
73) Anthracene-d10	17.430	188	1233536	37.127	ng/ul	0.00
81) Pyrene-d10	19.718	212	1466650	41.336	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.765	264	1379918	39.774	ng/ul	0.00

Target Compounds Qvalue

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

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