

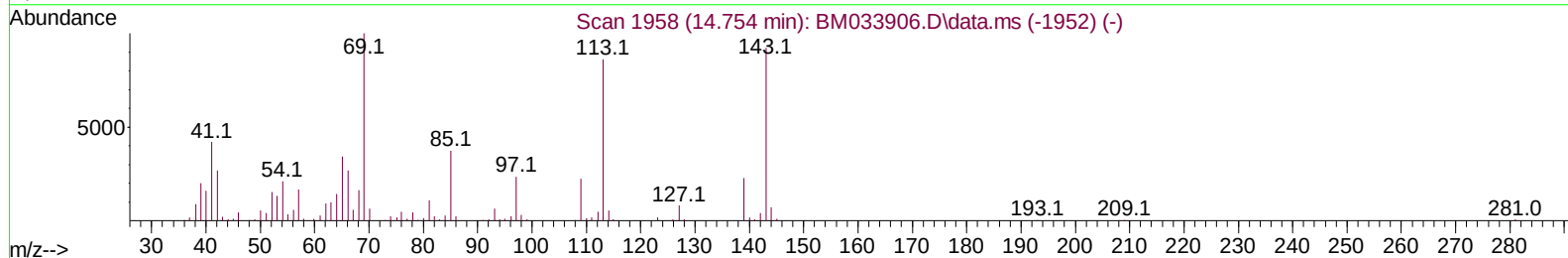
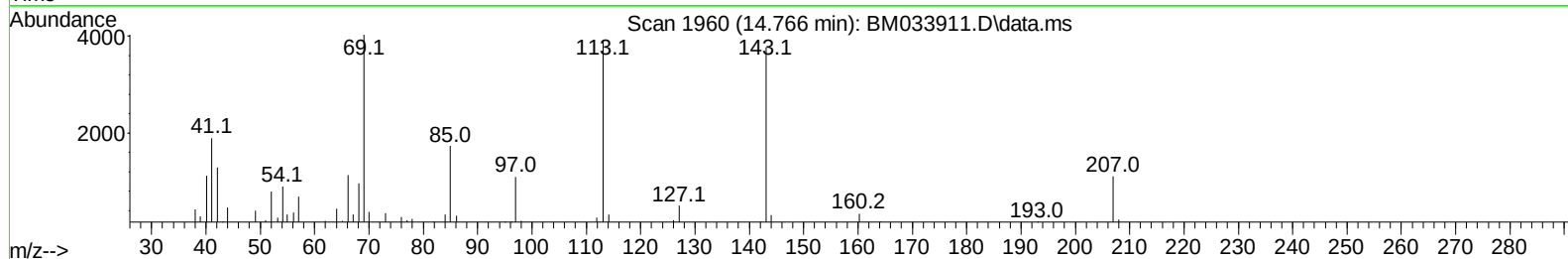
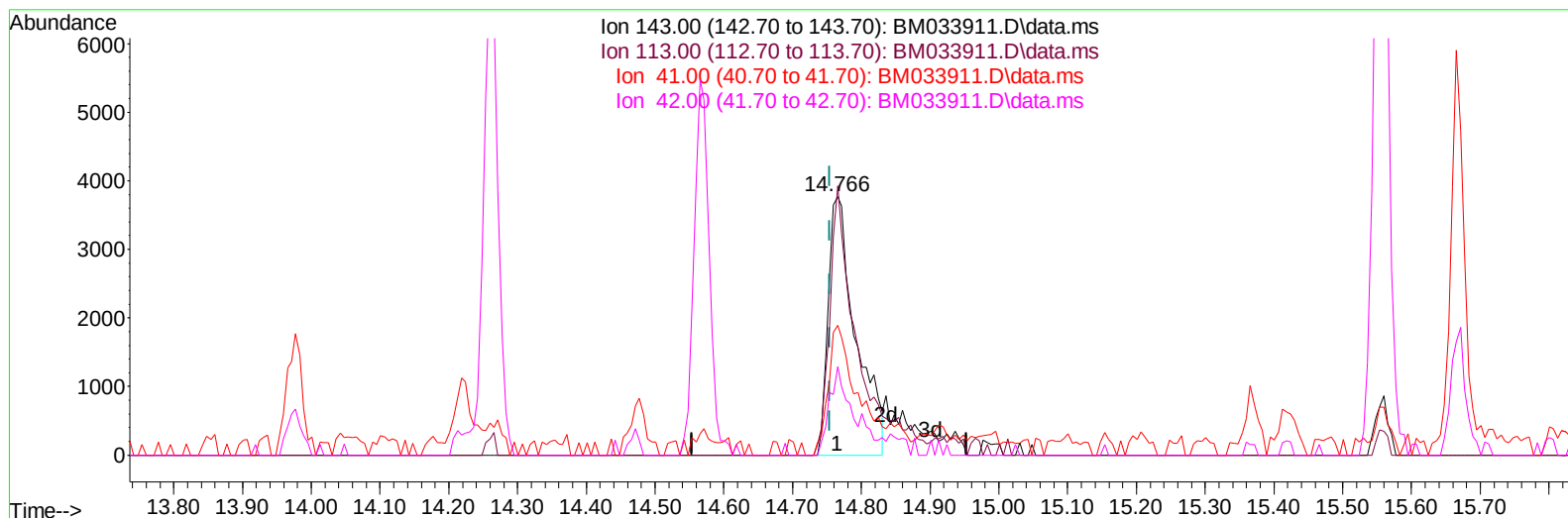
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
Data File : BM033911.D
Acq On : 28 Jan 2022 19:31
Operator : CG/JU
Sample : N1302-10
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0F29

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/31/2022
Supervised By :mohammad ahmed 02/01/2022

Quant Time: Jan 28 23:28:46 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 28 23:26:29 2022
Response via : Initial Calibration



TIC: BM033911.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 2.89 ng/ul

response 10511

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	104.09
41.00	57.20	50.27
42.00	39.50	34.16

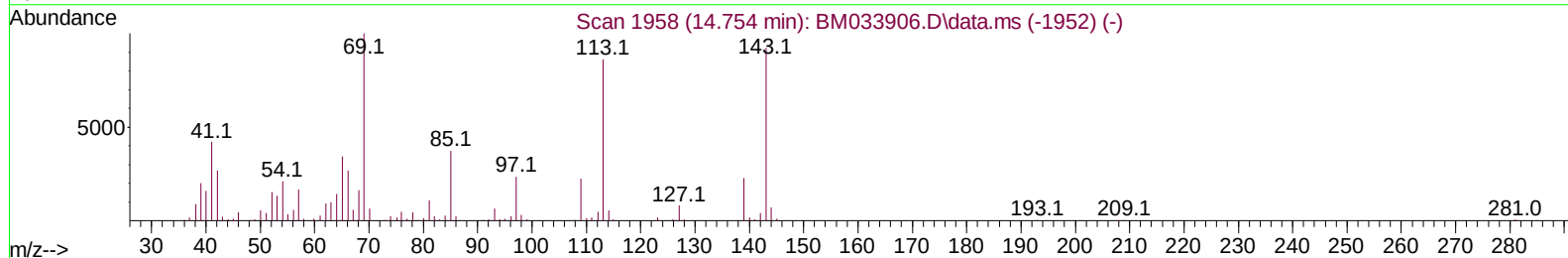
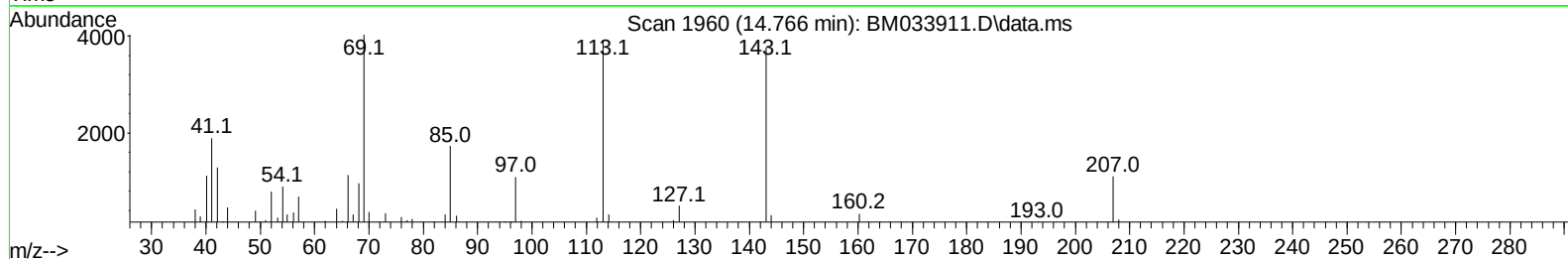
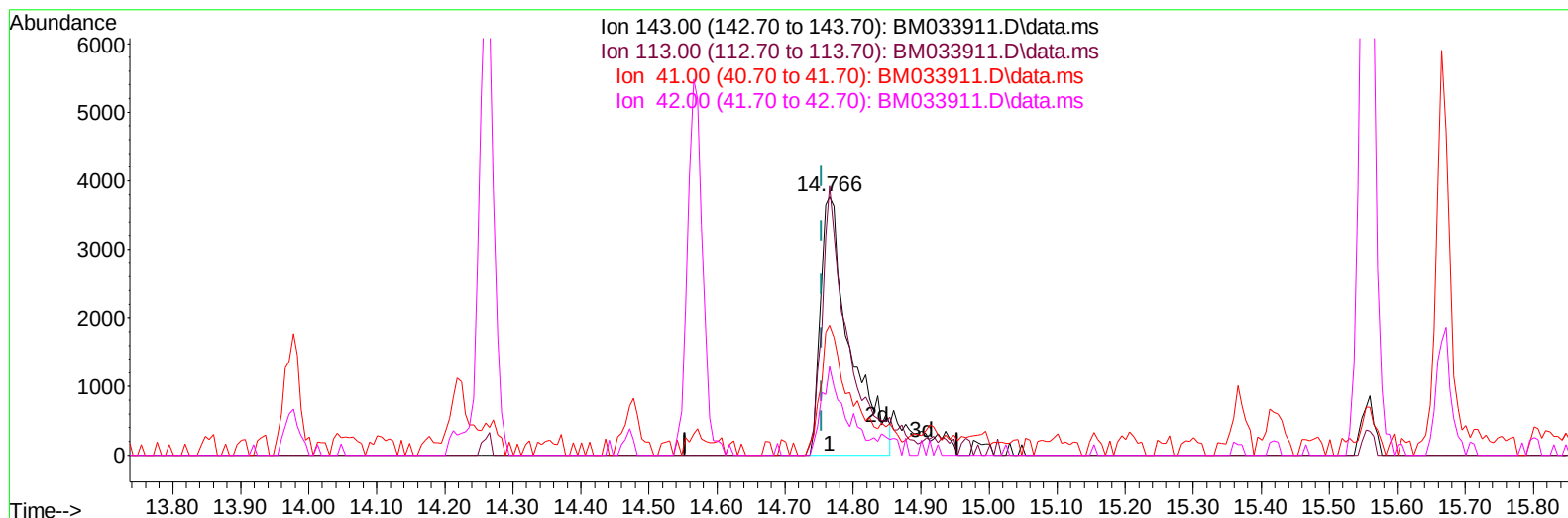
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TIC: BM033911.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.766min (+ 0.012) 3.12 ng/ul m

response 11338

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	104.09
41.00	57.20	50.27
42.00	39.50	34.16

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	98181	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	399821	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	240327	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.307	188	494990	20.000	ng/ul	0.00
79) Chrysene-d12	21.465	240	465087	20.000	ng/ul	# 0.00
88) Perylene-d12	23.859	264	476856	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	11345	4.341	ng/uL	0.00
4) Pyridine-d5	3.719	84	62325	8.588	ng/ul	0.00
7) Phenol-d5	7.084	99	50368	5.712	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	172060	32.041	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	167788	24.887	ng/ul	0.00
15) 4-Methylphenol-d8	8.637	113	98435	13.485	ng/ul	0.00
21) Nitrobenzene-d5	9.095	128	105648	33.459	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	107491	32.122	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	174188	27.152	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	254894	26.729	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	656636	34.679	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	768227	34.391	ng/ul	0.00
54) 4-Nitrophenol-d4	14.766	143	11338m	3.120	ng/ul	0.01
60) Fluorene-d10	15.560	176	560836	35.702	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.665	200	80275	27.264	ng/ul	0.00
73) Anthracene-d10	17.401	188	938062	39.055	ng/ul	0.00
81) Pyrene-d10	19.689	212	1067766	42.513	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.706	264	1022177	41.198	ng/ul	0.00

Target Compounds Qvalue

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

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