

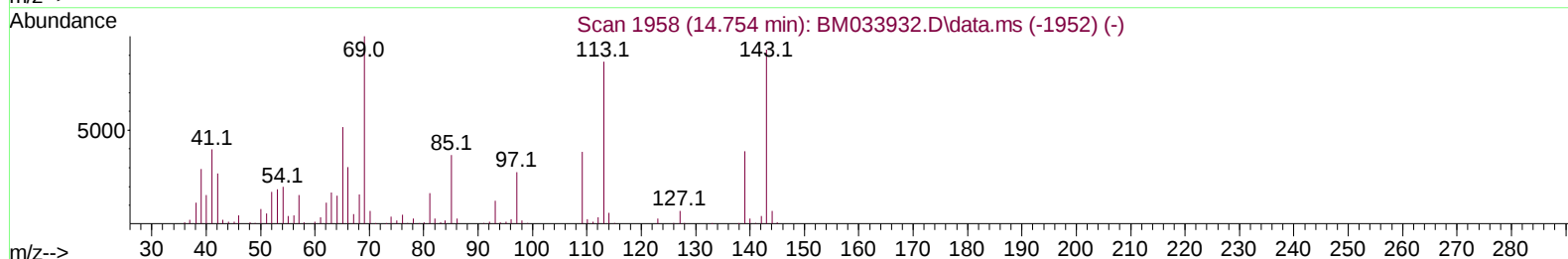
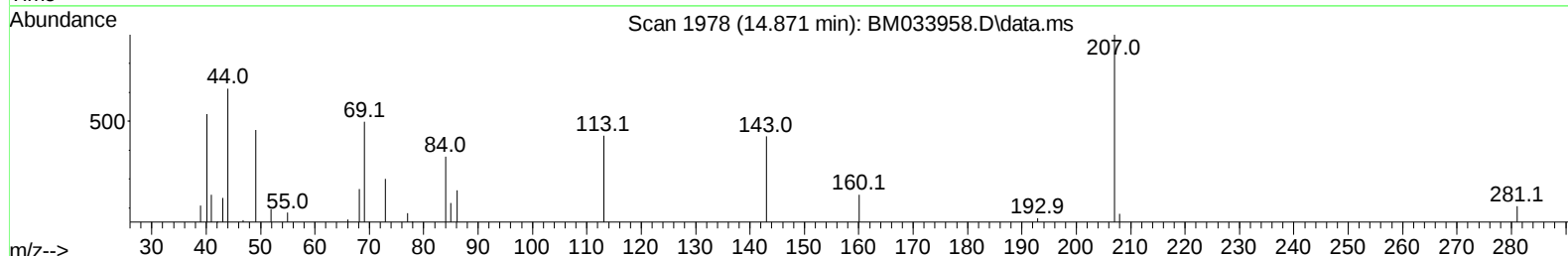
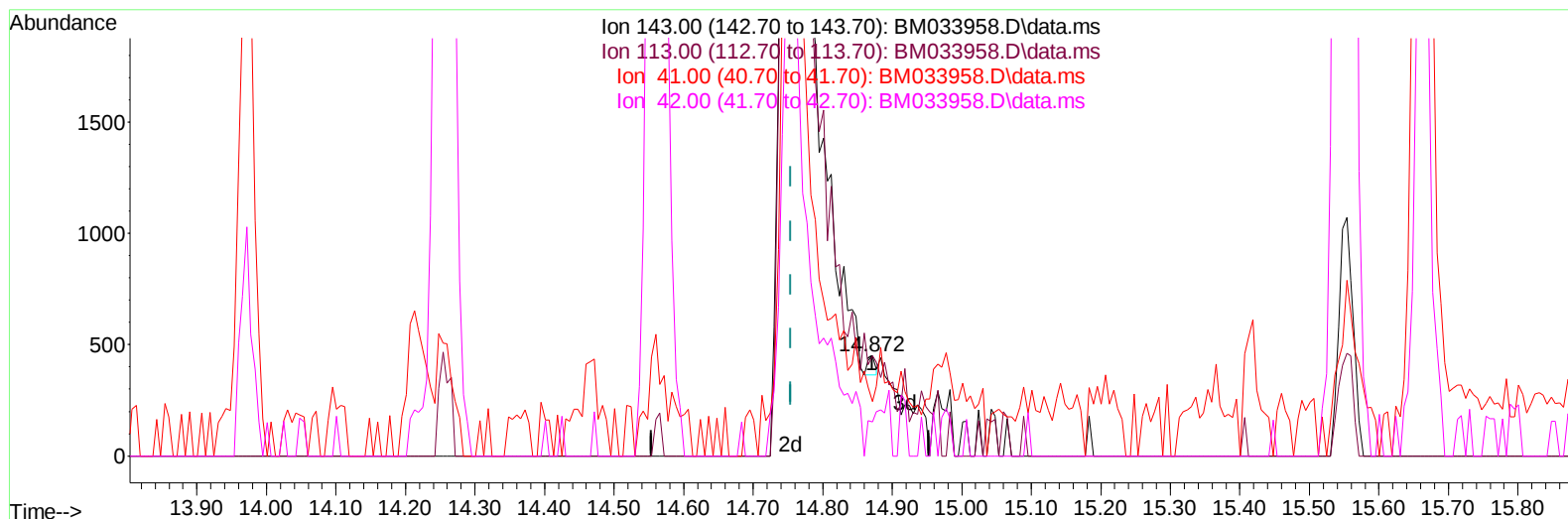
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033958.D
Acq On : 02 Feb 2022 02:49
Operator : CG/JU
Sample : N1307-02
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0Q11

Manual IntegrationsAPPROVED

Quant Time: Feb 02 04:08:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 01 02:16:28 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/02/2022
Supervised By :Yogesh Patel 02/04/2022



TIC: BM033958.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.871min (+ 0.118) 0.01 ng/u1

response 61

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	100.45
41.00	57.20	54.69
42.00	39.50	33.93

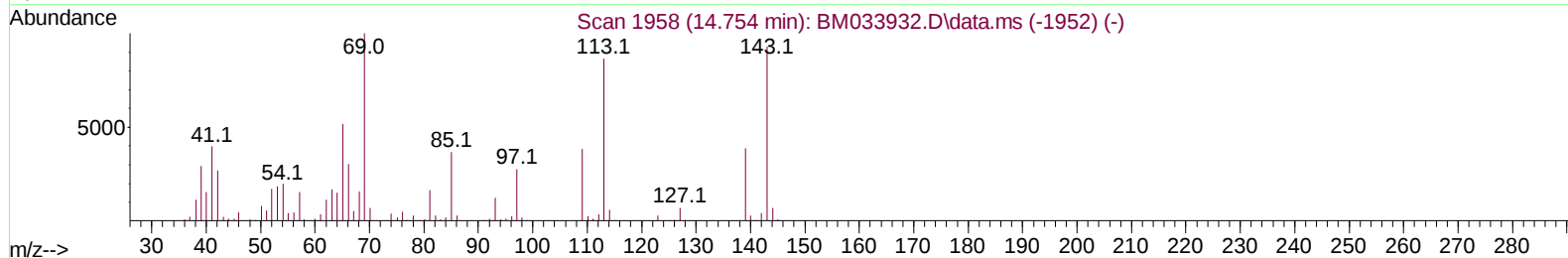
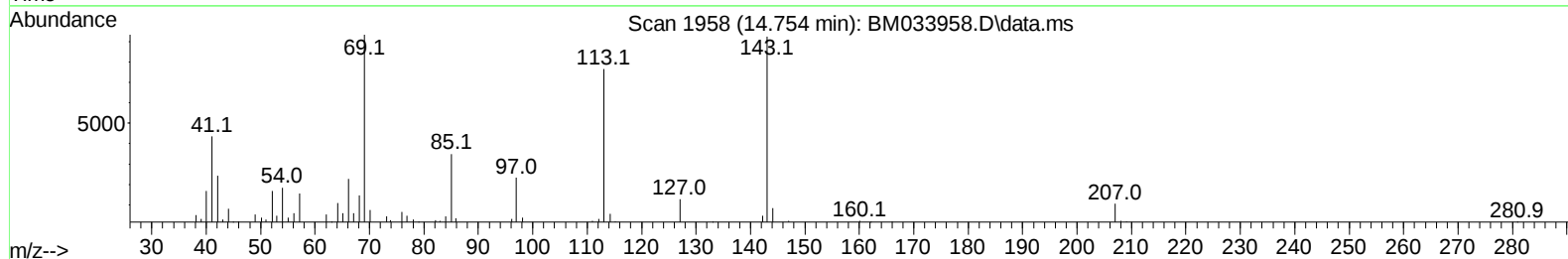
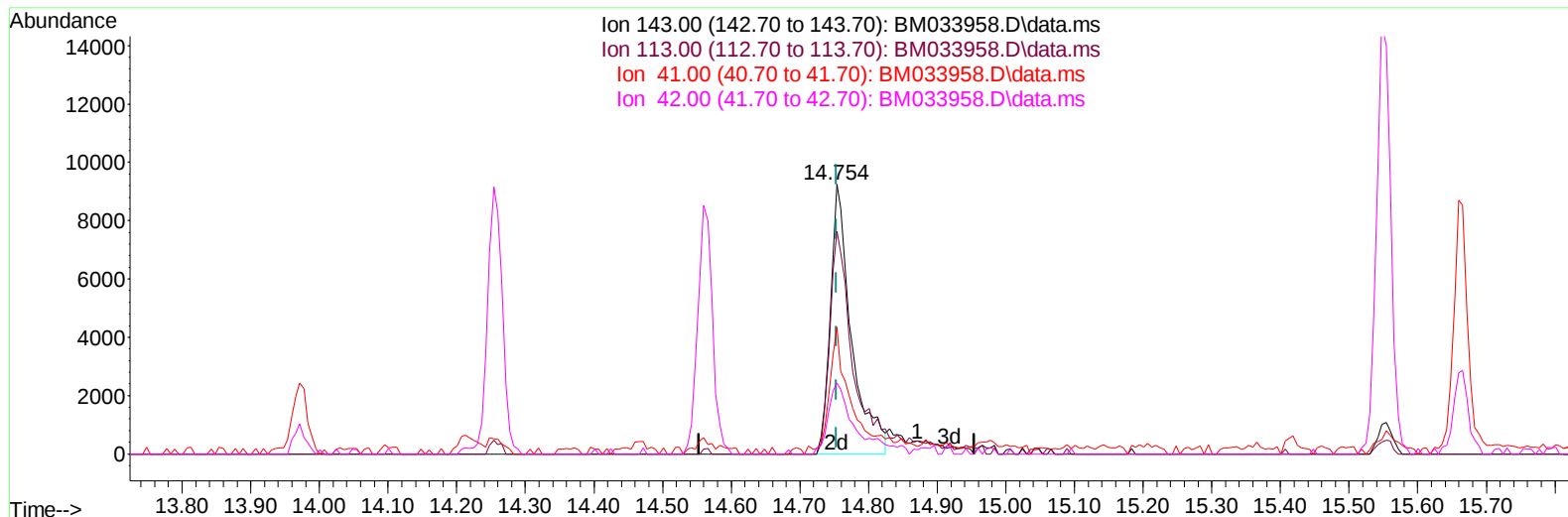
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(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.000) 4.30 ng/ul m

response 19775

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	82.73#
41.00	57.20	47.05
42.00	39.50	26.31#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	128290	20.000	ng/ul	0.00
20) Naphthalene-d8	10.736	136	558165	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	360041	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	686151	20.000	ng/ul	0.00
79) Chrysene-d12	21.459	240	534437	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	509431	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	14556	4.475	ng/uL	0.00
4) Pyridine-d5	3.731	84	29006	3.195	ng/ul	0.01
7) Phenol-d5	7.084	99	68999	6.420	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	215051	32.820	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	214295	24.887	ng/ul	0.00
15) 4-Methylphenol-d8	8.631	113	134738	15.821	ng/ul	-0.01
21) Nitrobenzene-d5	9.090	128	143190	31.971	ng/ul	0.00
24) 2-Nitrophenol-d4	9.813	143	152827	31.599	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165	239388	27.352	ng/ul	0.00
31) 4-Chloroaniline-d4	10.872	131	246977	20.435	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166	919898	34.191	ng/ul	0.00
49) Acenaphthylene-d8	14.260	160	1060192	32.167	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	19775m	4.302	ng/ul	0.00
60) Fluorene-d10	15.554	176	753765	33.552	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	130680	30.337	ng/ul	0.00
73) Anthracene-d10	17.401	188	1225394	37.292	ng/ul	0.00
81) Pyrene-d10	19.683	212	1270402	40.264	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.694	264	1031734	38.559	ng/ul	-0.01

Target Compounds Qvalue

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

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