

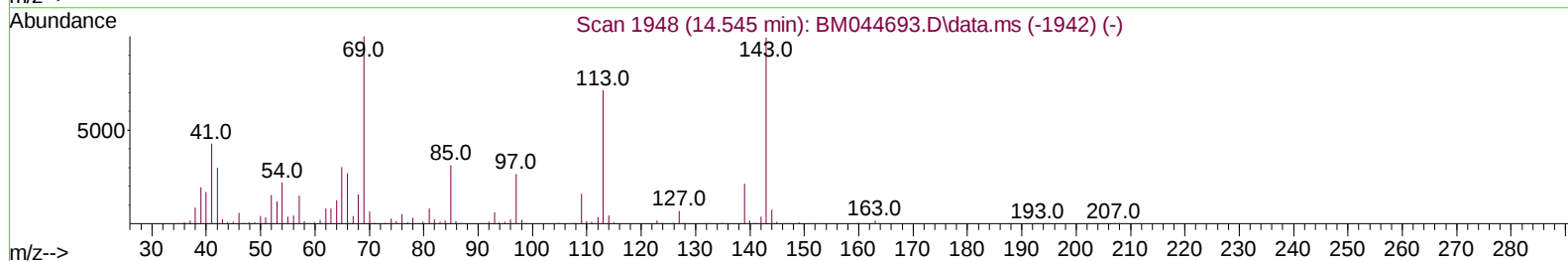
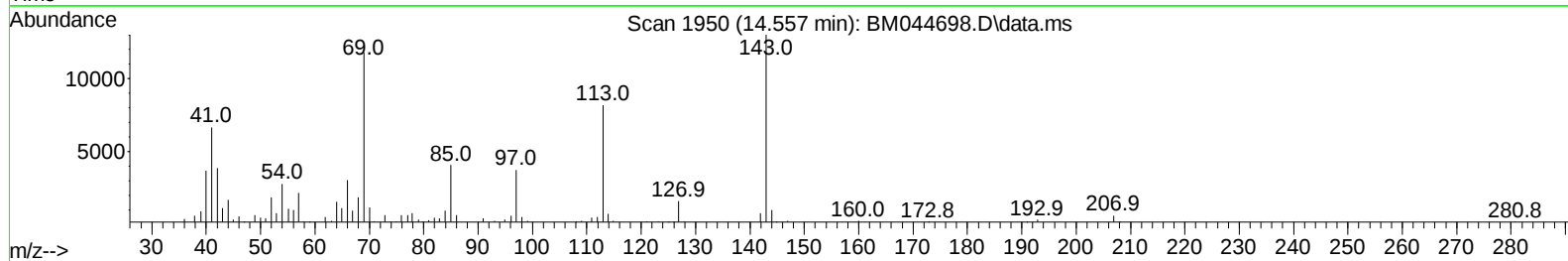
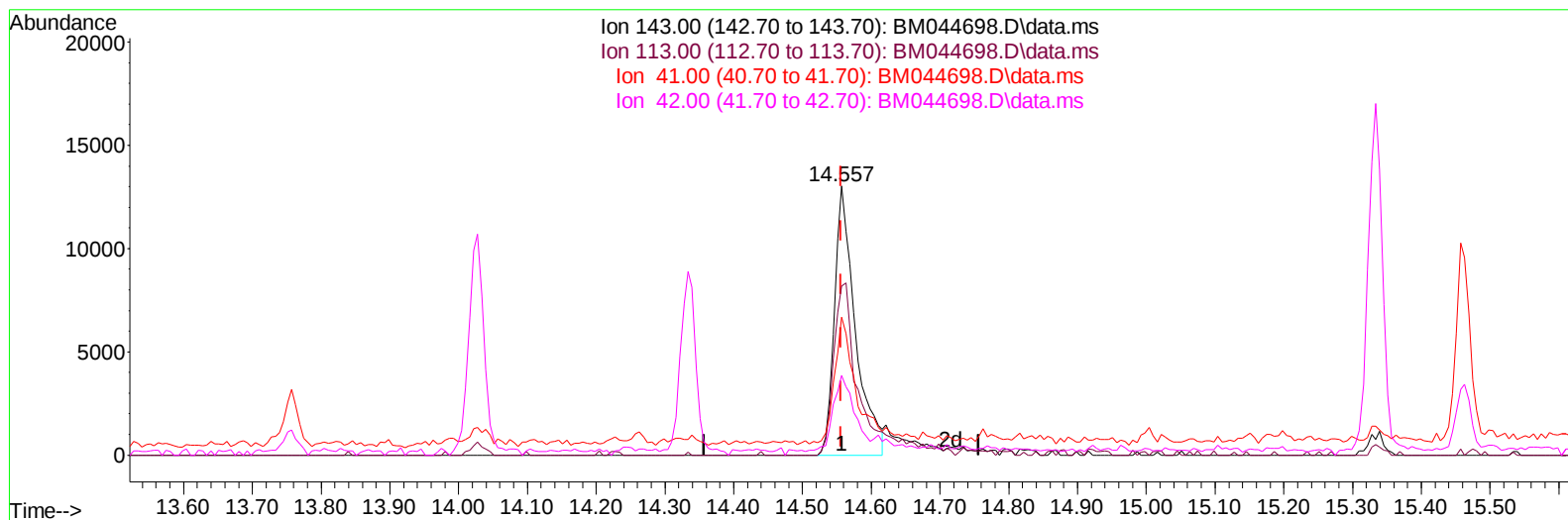
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032124\
Data File : BM044698.D
Acq On : 21 Mar 2024 13:14
Operator : MA/JU
Sample : P1813-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E0LX3

Manual IntegrationsAPPROVED

Quant Time: Mar 21 14:21:37 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 15 11:12:49 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/21/2024
Supervised By :mohammad ahmed 03/26/2024



TIC: BM044698.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.000) 4.82 ng/u1

response 27362

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	62.86
41.00	38.40	51.23#
42.00	26.60	29.60

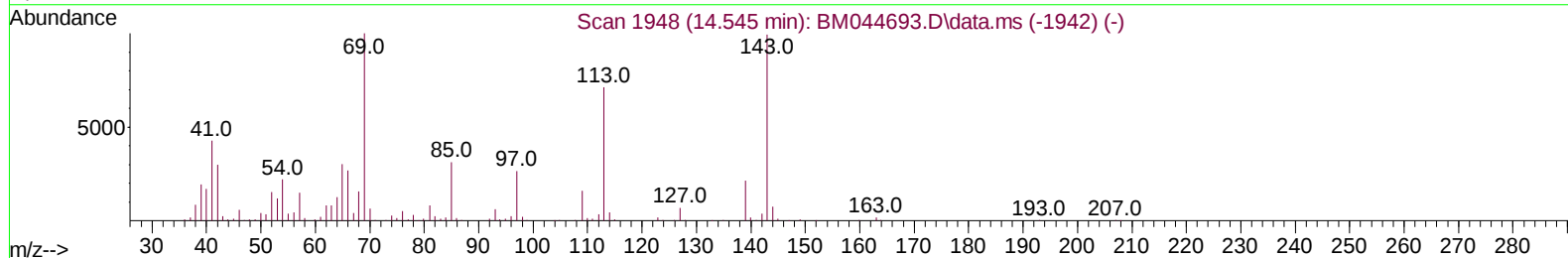
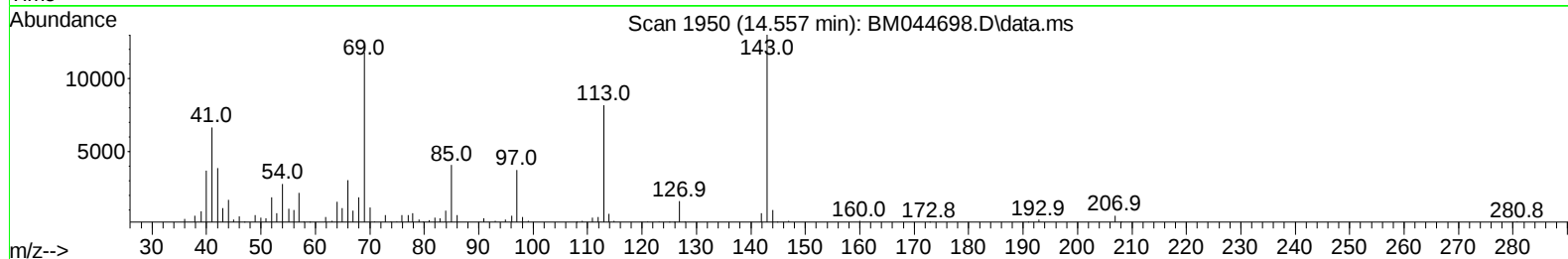
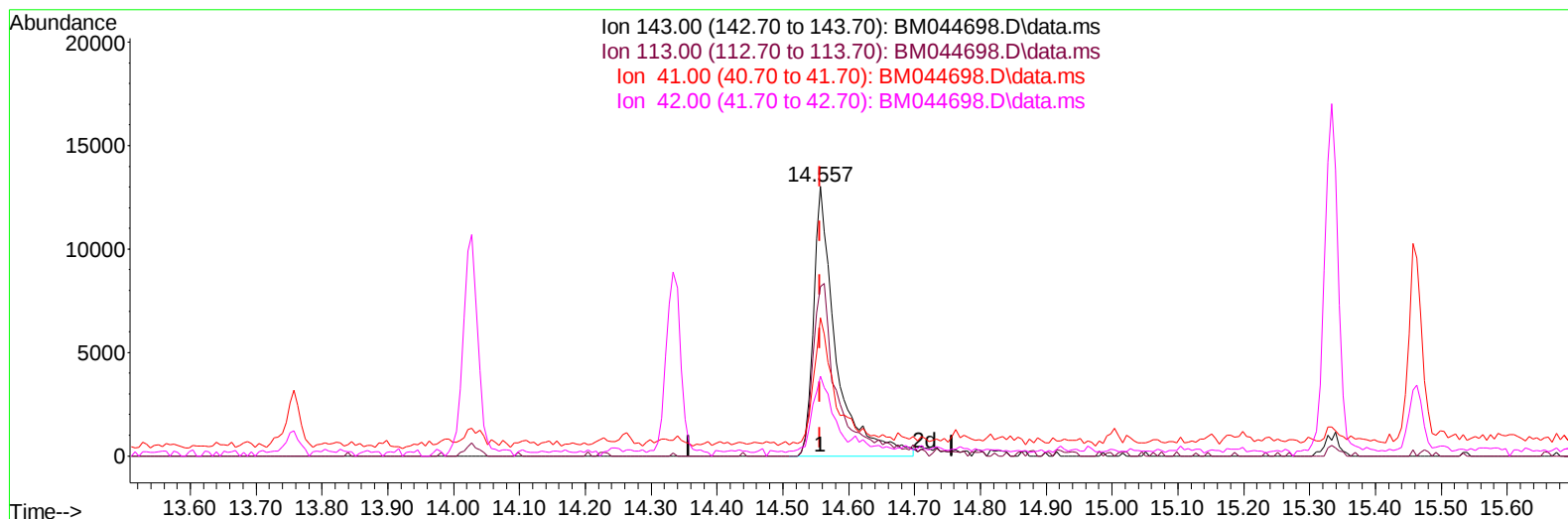
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TIC: BM044698.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.557min (+ 0.000) 5.43 ng/ul m

response 30814

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	63.70	62.86
41.00	38.40	51.23#
42.00	26.60	29.60

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Quant Time: Mar 21 14:22:19 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.693	152		158721	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136		630022	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.334	164		349125	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.092	188		648998	20.000	ng/ul	0.00
79) Chrysene-d12	21.286	240		538239	20.000	ng/ul	-0.01
88) Perylene-d12	23.533	264		642109	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.252	96		19811	4.233	ng/uL	0.00
4) Pyridine-d5	3.658	84		146609	10.644	ng/ul	0.00
7) Phenol-d5	6.863	99		118276	7.248	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67		294974	29.264	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.222	132		283280	23.623	ng/ul	-0.01
15) 4-Methylphenol-d8	8.393	113		214188	16.713	ng/ul	-0.01
21) Nitrobenzene-d5	8.852	128		169048	31.865	ng/ul	0.00
24) 2-Nitrophenol-d4	9.563	143		170044	29.732	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.093	165		290871	29.884	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.622	131		336026	20.786	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166		947510	34.598	ng/ul	-0.01
49) Acenaphthylene-d8	14.028	160		1071163	34.200	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143		30814m	5.426	ng/ul	0.00
60) Fluorene-d10	15.334	176		810484	35.216	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.463	200		128051	30.396	ng/ul	-0.01
73) Anthracene-d10	17.192	188		1212603	39.089	ng/ul	0.00
81) Pyrene-d10	19.492	212		1317106	39.898	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.386	264		1307967	38.755	ng/ul	-0.01

Target Compounds Qvalue

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

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