

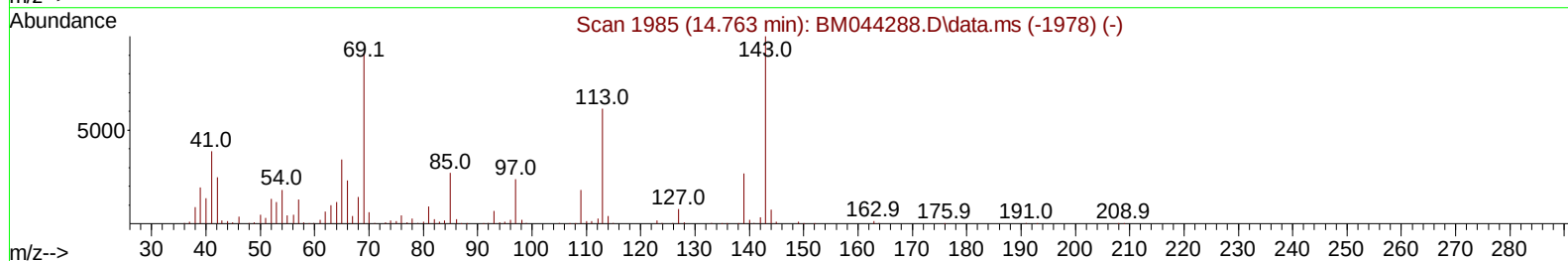
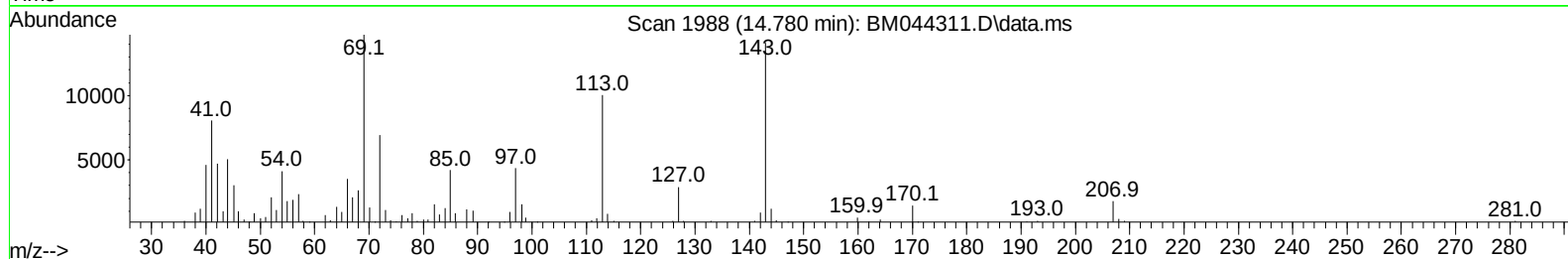
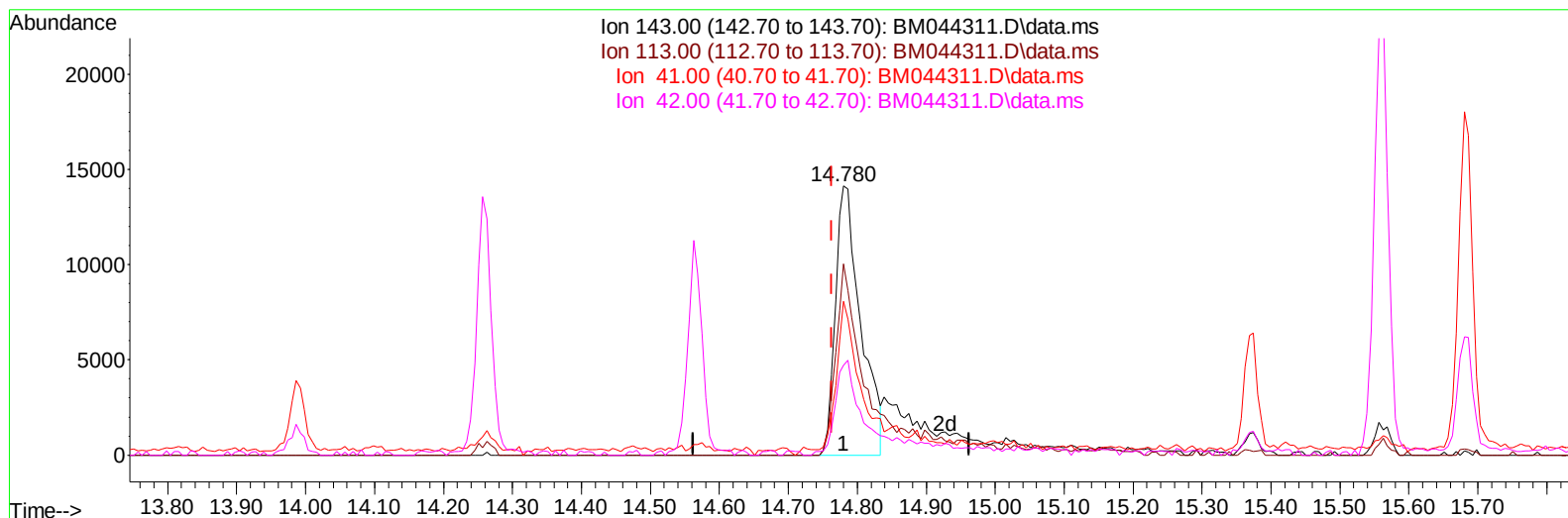
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022424\
Data File : BM044311.D
Acq On : 25 Feb 2024 01:22
Operator : MA/JU
Sample : P1494-13
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH3W3

Manual IntegrationsAPPROVED

Quant Time: Feb 25 23:05:02 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Feb 25 22:57:38 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 02/26/2024
Supervised By :mohammad ahmed 02/26/2024



TIC: BM044311.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.780min (+ 0.018) 4.49 ng/ul

response 35972

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	71.05#
41.00	38.20	57.09#
42.00	25.30	33.30#

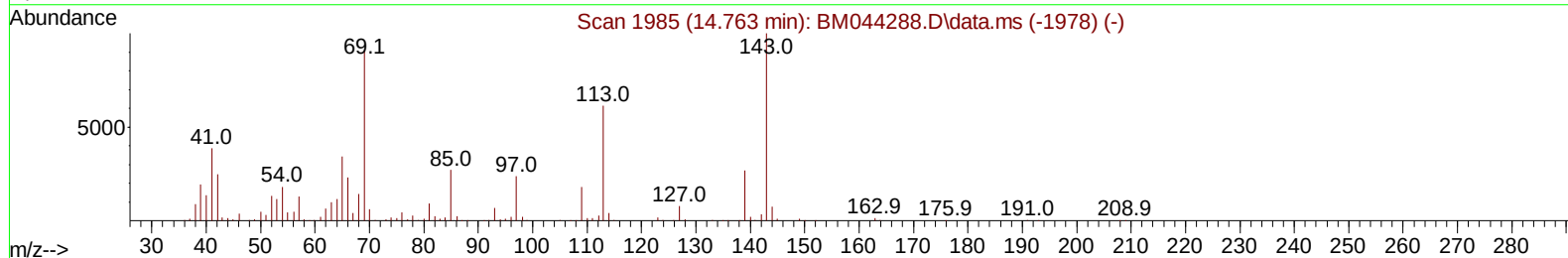
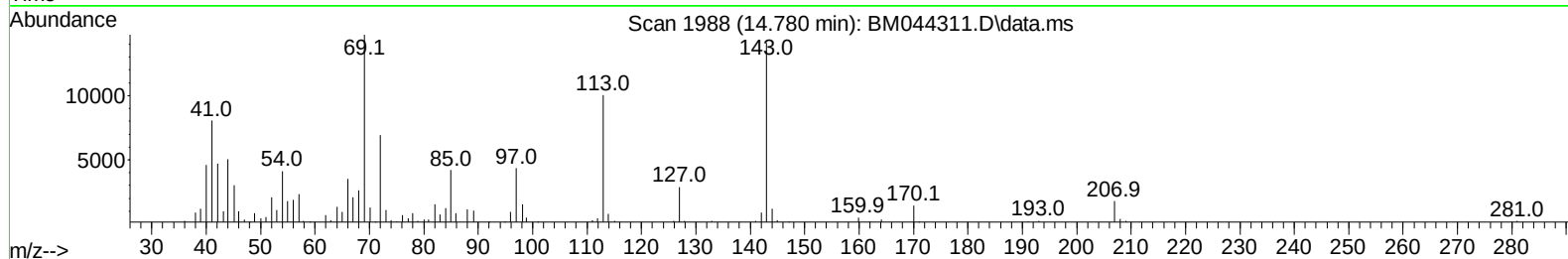
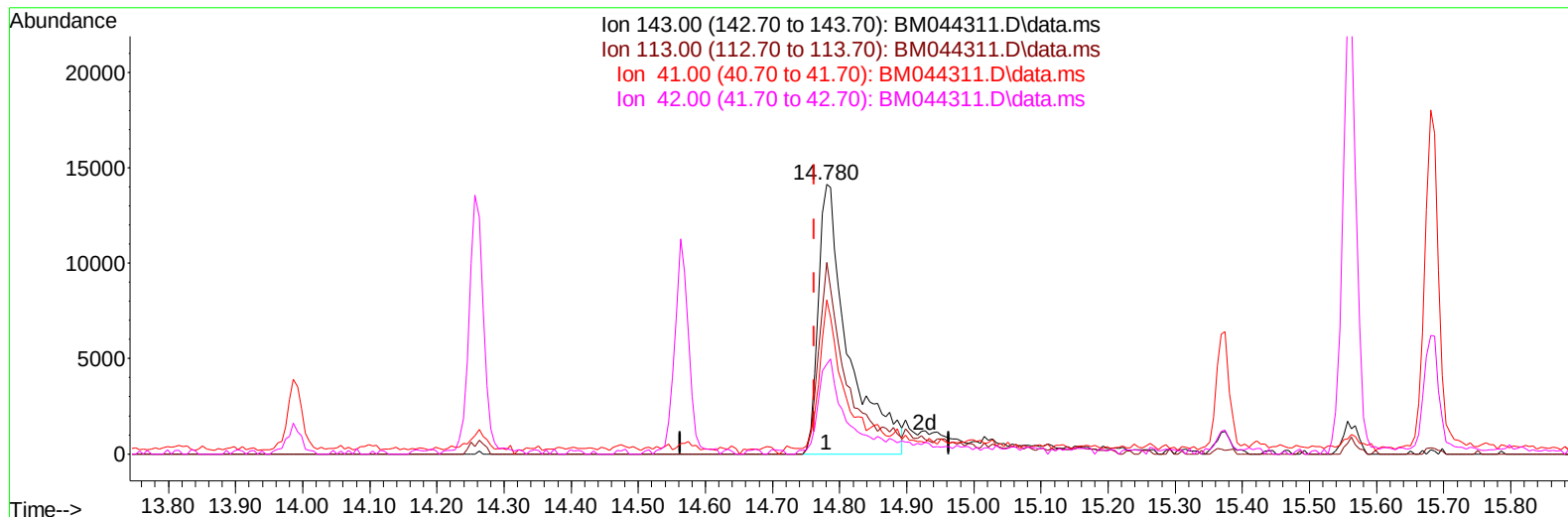
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TIC: BM044311.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.780min (+ 0.018) 5.46 ng/ul m

response 43717

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	71.05#
41.00	38.20	57.09#
42.00	25.30	33.30#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 02/26/2024
 Supervised By :mohammad ahmed 02/26/2024

Quant Time: Feb 26 01:08:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Feb 25 22:57:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	203588	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	886336	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.563	164	522042	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	1121220	20.000	ng/ul	0.00
79) Chrysene-d12	21.515	240	1008111	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	1135080	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	25634	3.973	ng/uL	0.00
4) Pyridine-d5	3.775	84	189859	10.774	ng/ul	0.00
7) Phenol-d5	7.087	99	133502	6.845	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	436787	34.585	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	397546	27.065	ng/ul	0.01
15) 4-Methylphenol-d8	8.634	113	242578	16.438	ng/ul	0.00
21) Nitrobenzene-d5	9.092	128	259714	36.038	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	262658	35.724	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	407171	31.394	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	612818	28.133	ng/ul	0.01
46) Dimethylphthalate-d6	13.992	166	1544641	38.365	ng/ul	0.01
49) Acenaphthylene-d8	14.257	160	1763068	38.060	ng/ul	0.00
54) 4-Nitrophenol-d4	14.780	143	43717m	5.457	ng/ul	0.02
60) Fluorene-d10	15.563	176	1388857	40.755	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	258200	37.114	ng/ul	0.00
73) Anthracene-d10	17.415	188	2465566	46.119	ng/ul	0.00
81) Pyrene-d10	19.715	212	2865331	47.244	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	2766779	47.570	ng/ul	0.00

Target Compounds Qvalue

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

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