

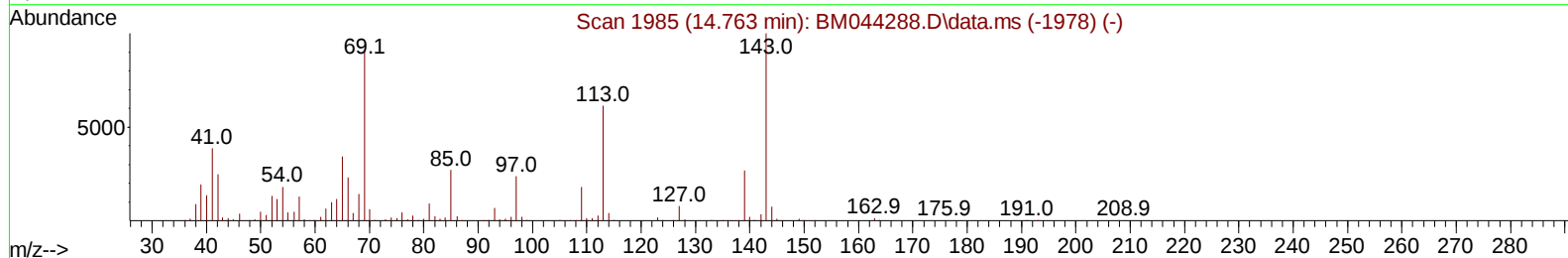
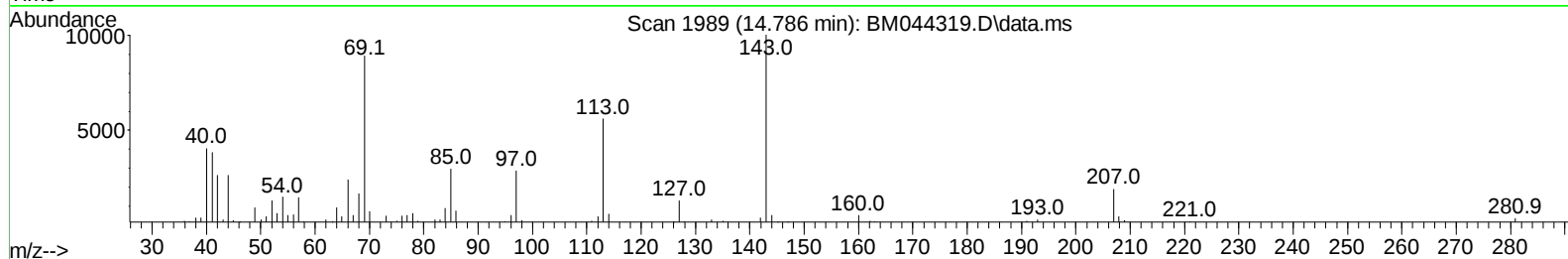
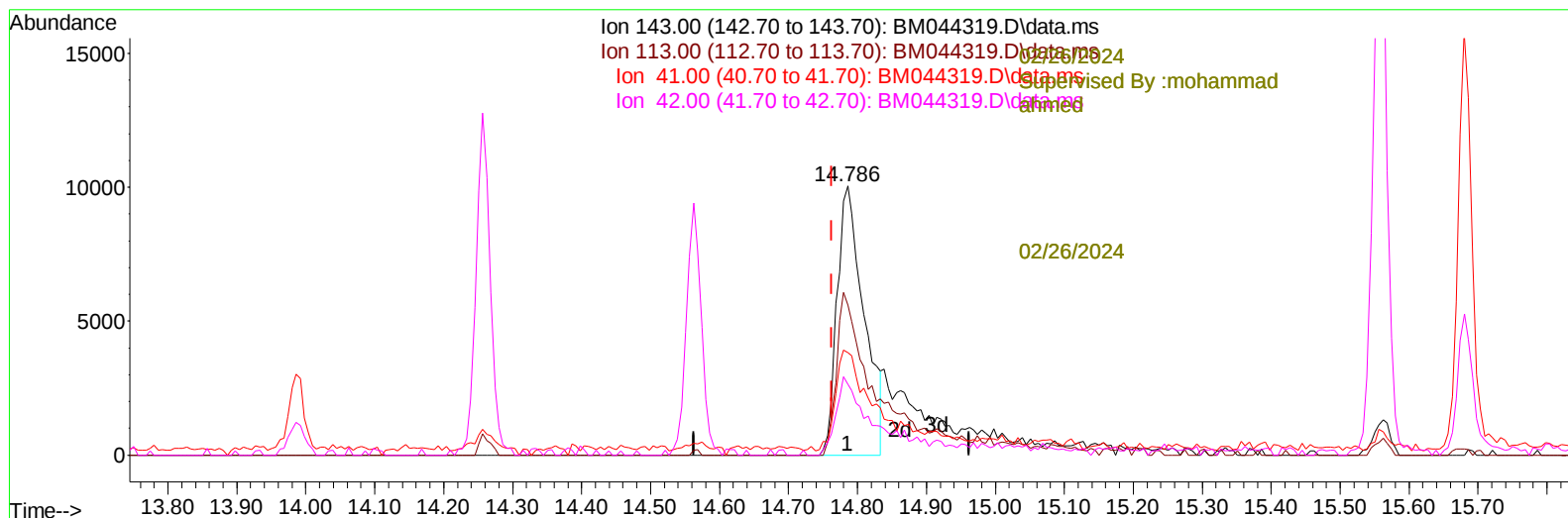
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022424\
Data File : BM044319.D
Acq On : 25 Feb 2024 06:12
Operator : MA/JU
Sample : P1494-10
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH3W0

Manual IntegrationsAPPROVED

Quant Time: Feb 25 23:07:11 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: BM044319.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.786min (+ 0.023) 3.82 ng/u1

response 27220

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	55.79
41.00	38.20	38.16
42.00	25.30	26.20

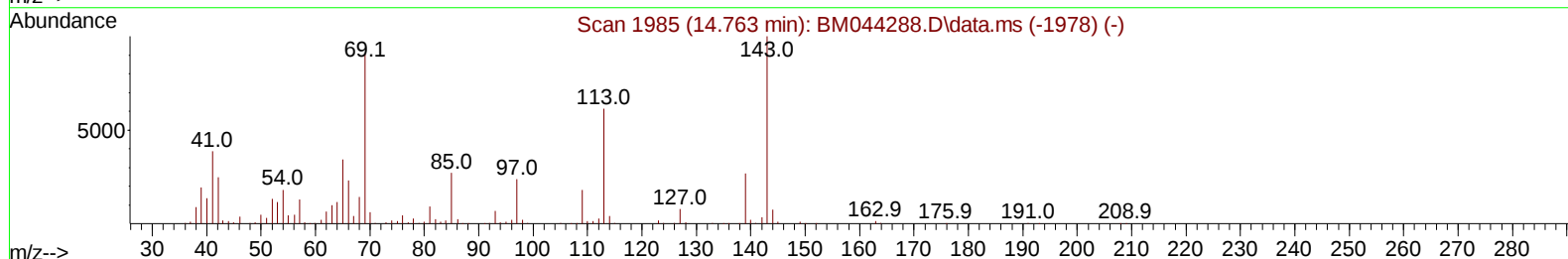
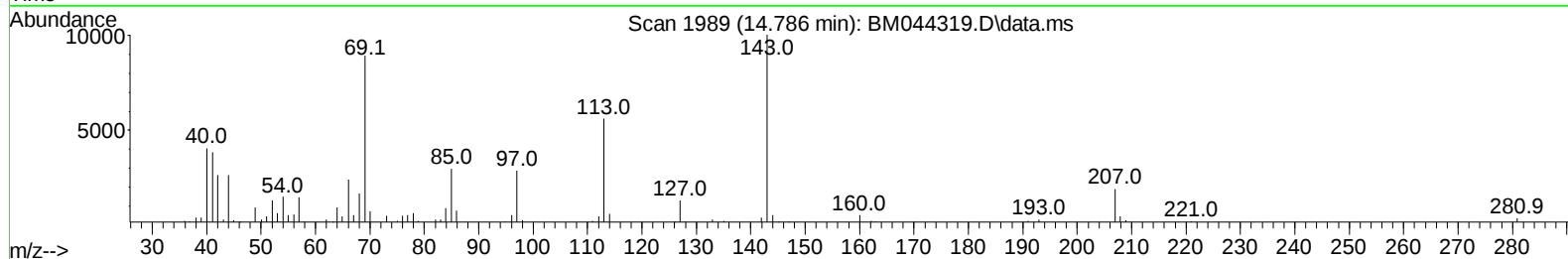
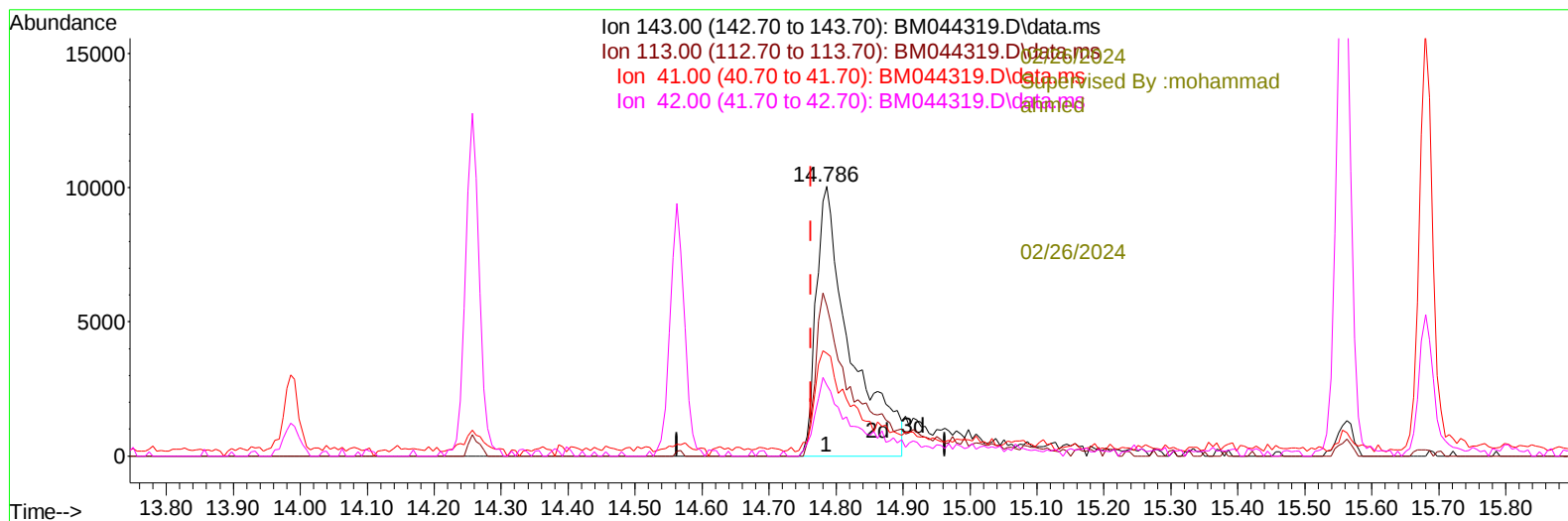
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TIC: BM044319.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.786min (+ 0.023) 4.98 ng/ul m

response 35427

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	55.79
41.00	38.20	38.16
42.00	25.30	26.20

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Manual IntegrationsAPPROVED

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

Quant Time: Feb 26 01:34:34 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)
-----02/26/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	190126	20.000	ng/ul	0.00
20) Naphthalene-d8	10.727	136	803724	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.562	164	463931	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	1002067	20.000	ng/ul	0.00
79) Chrysene-d12	21.515	240	909013	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	1042289	20.000	ng/ul	0.00
-----02/26/2024						
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	21403	3.552	ng/uL	0.00
4) Pyridine-d5	3.781	84	113457	6.894	ng/ul	0.01
7) Phenol-d5	7.086	99	122404	6.721	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	391908	33.228	ng/ul	0.00
11) 2-Chlorophenol-d4	7.445	132	368187	26.841	ng/ul	0.00
15) 4-Methylphenol-d8	8.633	113	215304	15.623	ng/ul	0.00
21) Nitrobenzene-d5	9.092	128	228131	34.910	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	230974	34.643	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.345	165	348515	29.633	ng/ul	0.00
31) 4-Chloroaniline-d4	10.874	131	517158	26.182	ng/ul	0.01
46) Dimethylphthalate-d6	13.986	166	1318134	36.840	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1523581	37.010	ng/ul	0.00
54) 4-Nitrophenol-d4	14.786	143	35427m	4.976	ng/ul	0.02
60) Fluorene-d10	15.557	176	1223255	40.392	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.680	200	224872	36.167	ng/ul	0.00
73) Anthracene-d10	17.415	188	2226625	46.602	ng/ul	0.00
81) Pyrene-d10	19.715	212	2543743	46.514	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.768	264	2521403	47.210	ng/ul	0.00

Target Compounds Qvalue

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

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