

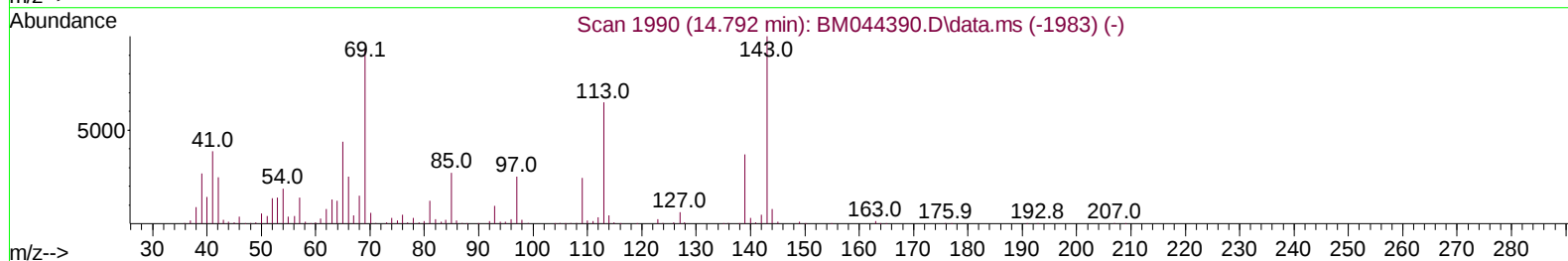
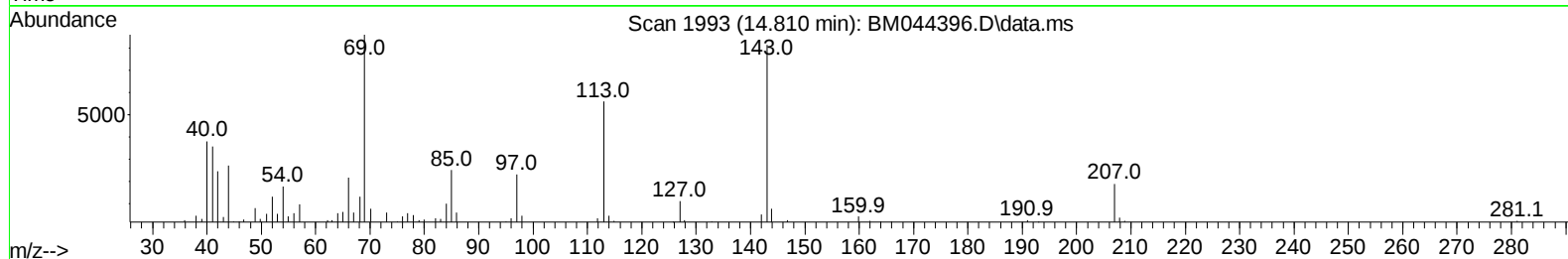
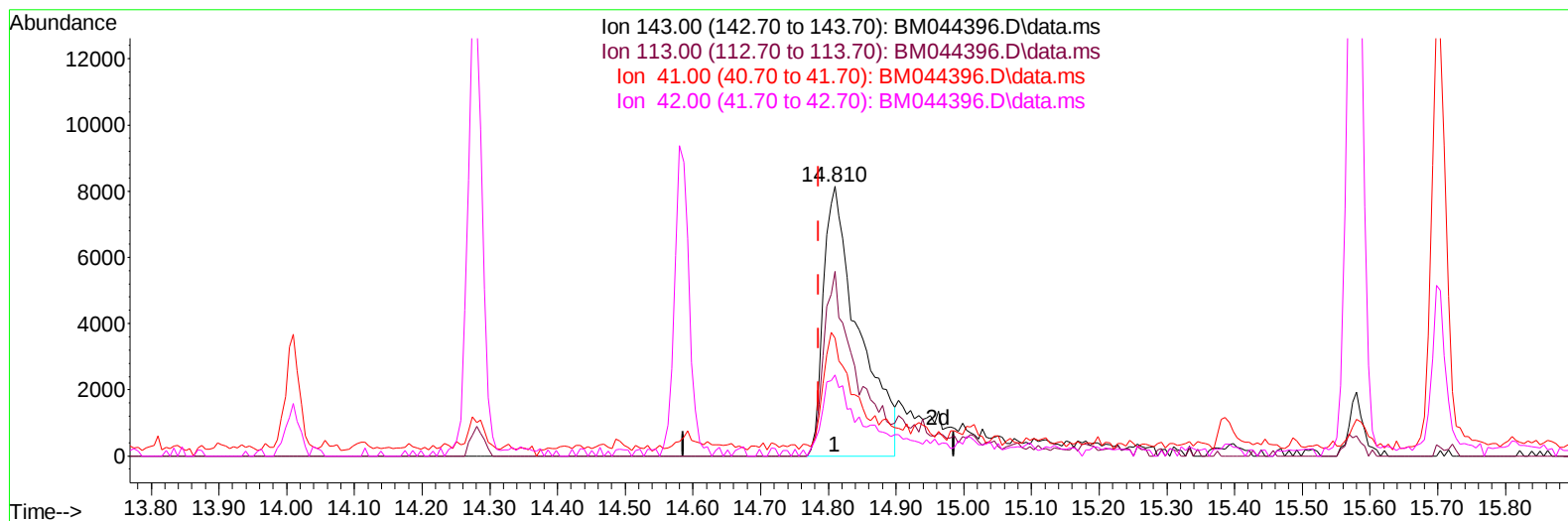
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
Data File : BM044396.D
Acq On : 28 Feb 2024 10:09
Operator : MA/JU
Sample : P1508-09
Misc :
ALS Vial : 78 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH409

Manual IntegrationsAPPROVED

Quant Time: Feb 28 11:40:11 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 27 23:05:48 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/28/2024
Supervised By :mohammad ahmed 02/29/2024



TIC: BM044396.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.024) 4.06 ng/u1

response 29303

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	68.52
41.00	38.20	43.86
42.00	25.30	30.09

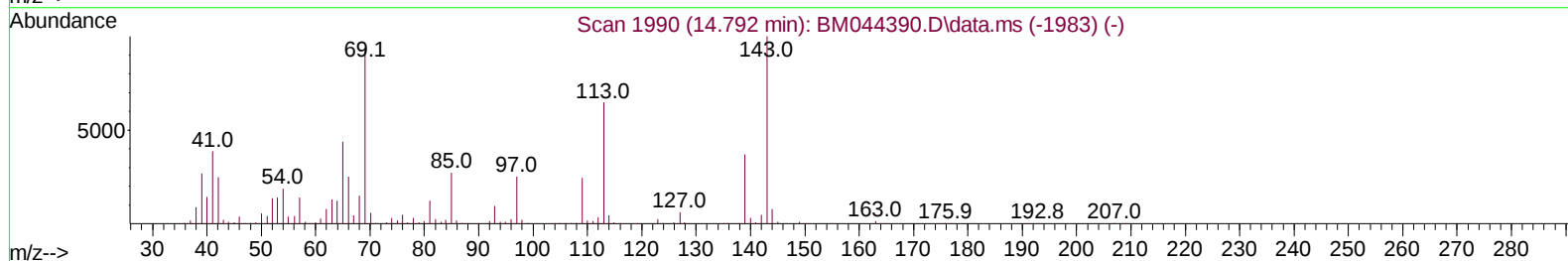
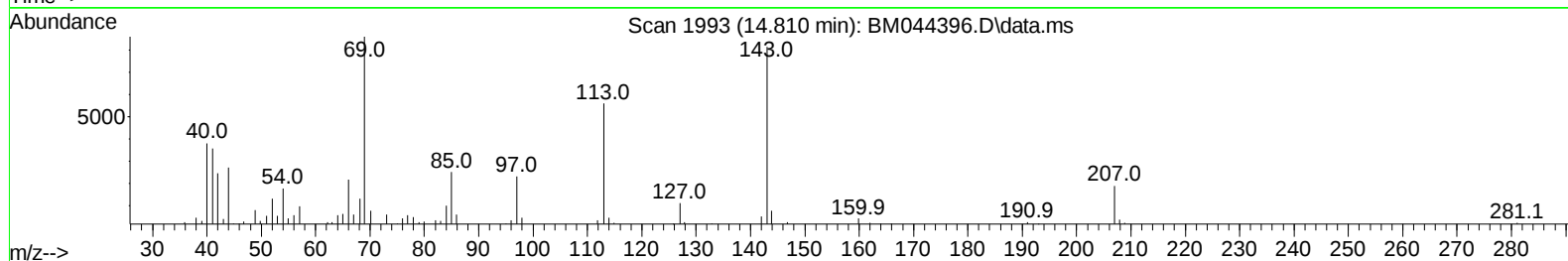
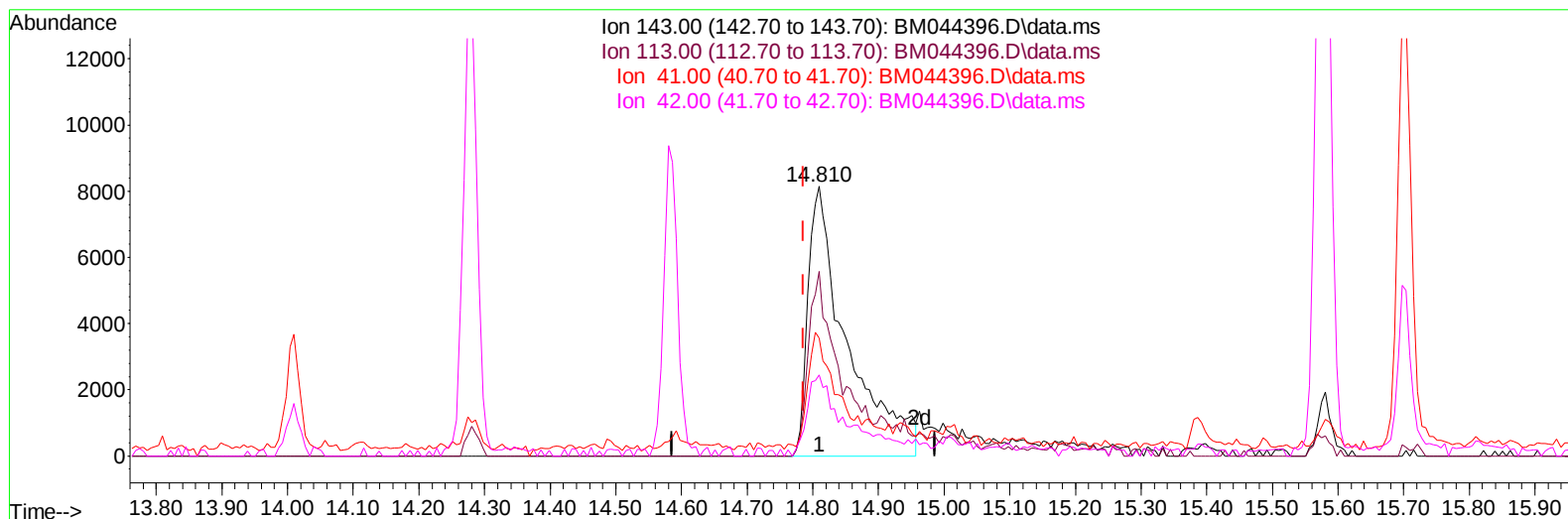
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TIC: BM044396.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.024) 4.68 ng/ul m

response 33755

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	68.52
41.00	38.20	43.86
42.00	25.30	30.09

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	186594	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	810935	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	470100	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	952239	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	757553	20.000	ng/ul	0.00
88) Perylene-d12	23.950	264	845927	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	27374	4.629	ng/uL	0.00
4) Pyridine-d5	3.793	84	148761	9.210	ng/ul	0.01
7) Phenol-d5	7.110	99	135460	7.578	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	449220	38.808	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	417907	31.042	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	248810	18.396	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	274435	41.622	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	280105	41.639	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	431553	36.368	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	608039	30.509	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1541926	42.529	ng/ul	0.00
49) Acenaphthylene-d8	14.280	160	1805546	43.284	ng/ul	0.00
54) 4-Nitrophenol-d4	14.810	143	33755m	4.679	ng/ul	0.02
60) Fluorene-d10	15.580	176	1398210	45.563	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	212122	35.901	ng/ul	0.00
73) Anthracene-d10	17.433	188	2290701	50.451	ng/ul	0.00
81) Pyrene-d10	19.733	212	2490910	54.655	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.797	264	2220367	51.224	ng/ul	0.00

Target Compounds Qvalue

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

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