

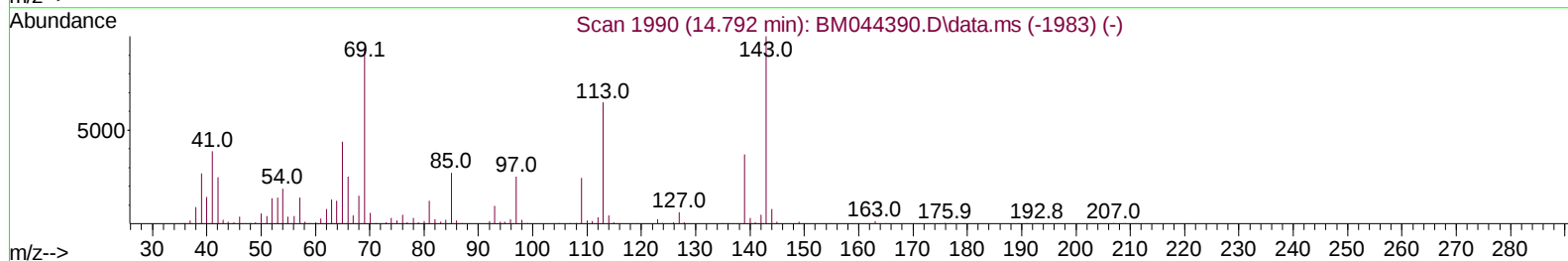
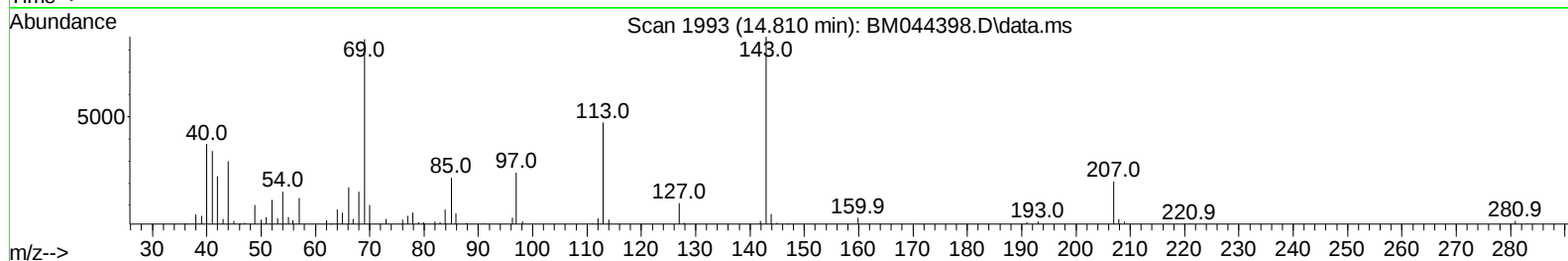
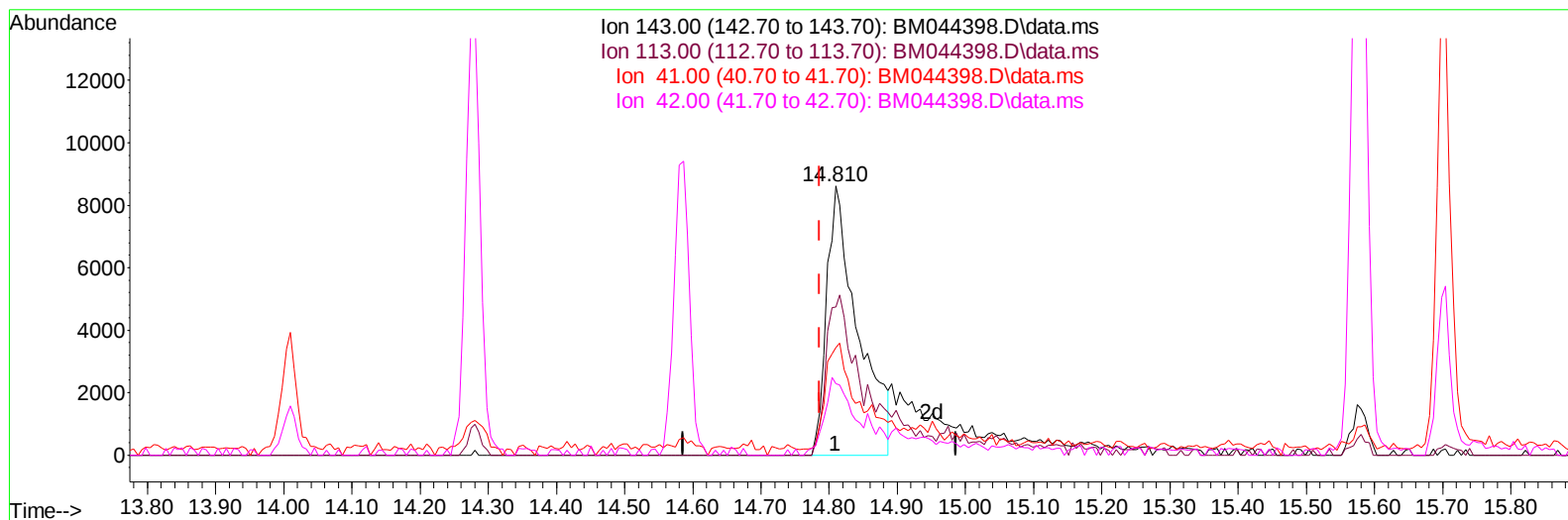
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022624\
 Data File : BM044398.D
 Acq On : 28 Feb 2024 11:22
 Operator : MA/JU
 Sample : P1508-01
 Misc :
 ALS Vial : 80 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH3Z5

Manual IntegrationsAPPROVED

Quant Time: Feb 28 12:17:06 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: BM044398.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.024) 3.46 ng/u1

response 27413

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	55.00
41.00	38.20	40.01
42.00	25.30	26.65

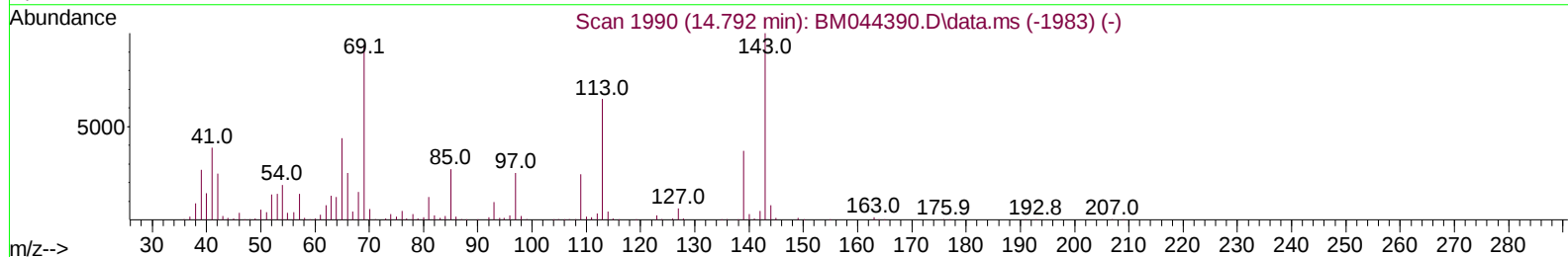
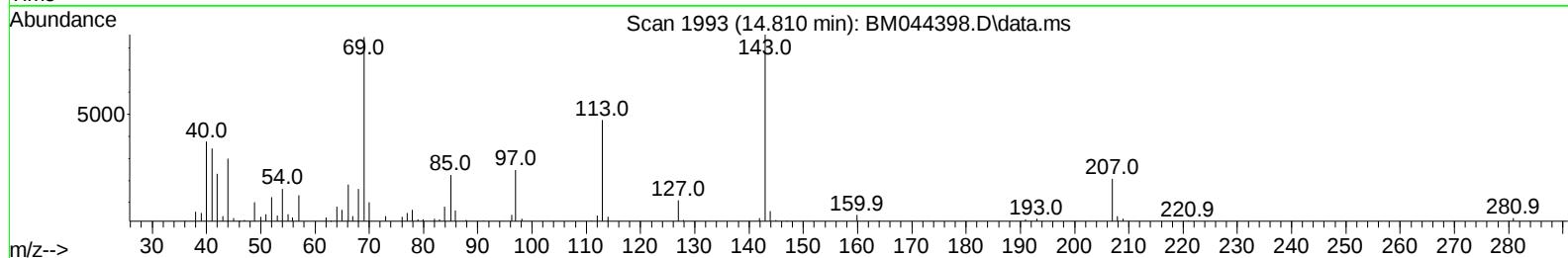
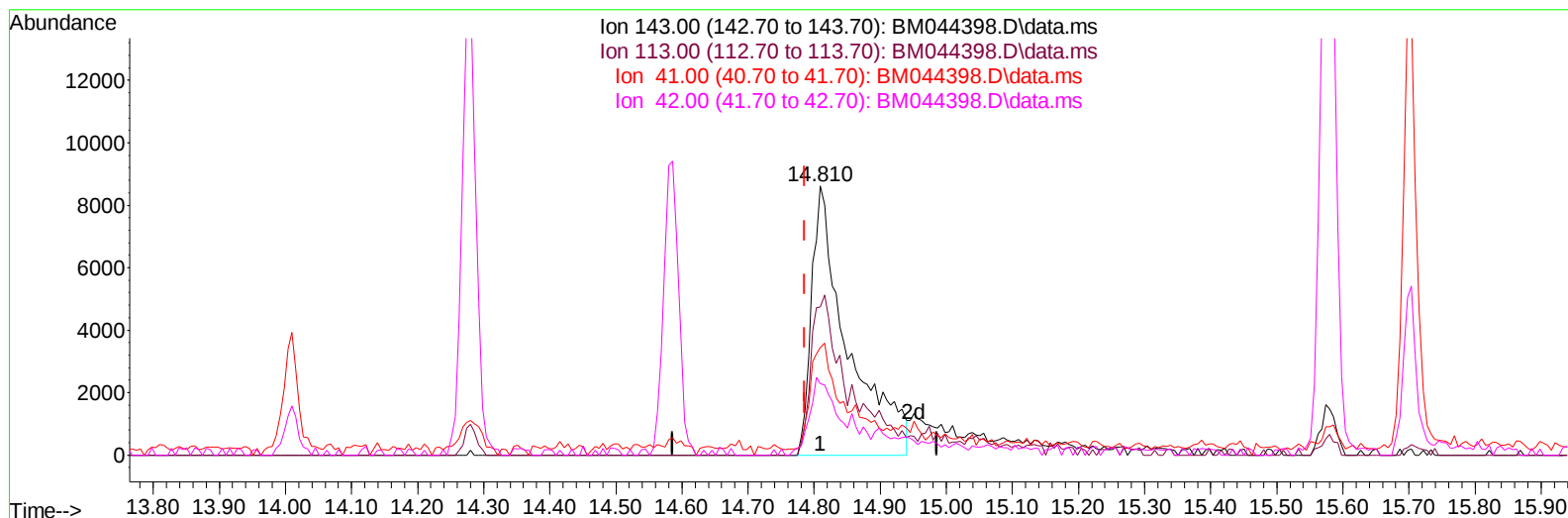
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(54) 4-Nitrophenol-d4 (S)

14.810min (+ 0.024) 4.13 ng/ul m

response 32684

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	55.00
41.00	38.20	40.01
42.00	25.30	26.65

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Reviewed By :Jagrut Upadhyay 02/28/2024
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Quant Time: Feb 28 12:18:03 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	197000	20.000	ng/ul	0.00
20) Naphthalene-d8	10.751	136	871797	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	515996	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.333	188	1059535	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	872246	20.000	ng/ul	0.00
88) Perylene-d12	23.950	264	980297	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	27257	4.365	ng/uL	0.00
4) Pyridine-d5	3.787	84	205282	12.039	ng/ul	0.00
7) Phenol-d5	7.110	99	136558	7.236	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	458923	37.552	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	425763	29.955	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	255156	17.869	ng/ul	0.00
21) Nitrobenzene-d5	9.116	128	283272	39.963	ng/ul	0.00
24) 2-Nitrophenol-d4	9.833	143	289450	40.024	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	449649	35.247	ng/ul	0.00
31) 4-Chloroaniline-d4	10.898	131	658097	30.716	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166	1578041	39.653	ng/ul	0.00
49) Acenaphthylene-d8	14.280	160	1869352	40.827	ng/ul	0.00
54) 4-Nitrophenol-d4	14.810	143	32684m	4.127	ng/ul	0.02
60) Fluorene-d10	15.580	176	1443383	42.852	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	215129	32.723	ng/ul	0.00
73) Anthracene-d10	17.433	188	2409737	47.698	ng/ul	0.00
81) Pyrene-d10	19.733	212	2640962	50.327	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.797	264	2449770	48.770	ng/ul	0.00

Target Compounds Qvalue

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

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