

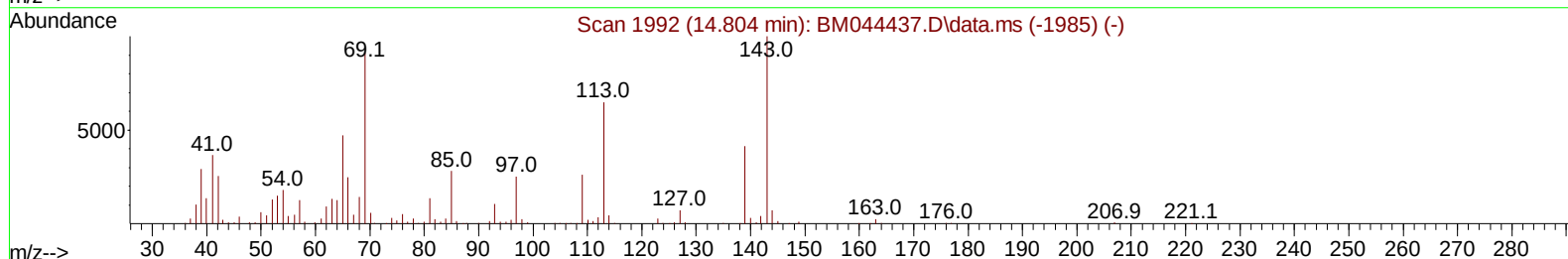
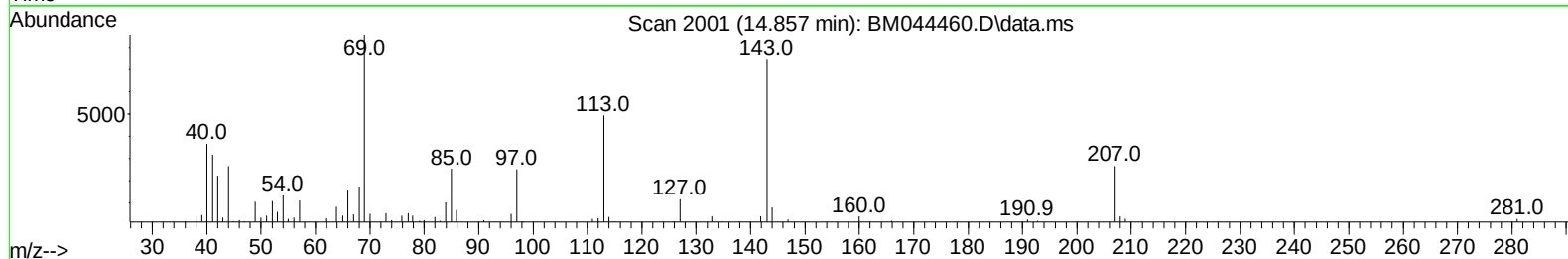
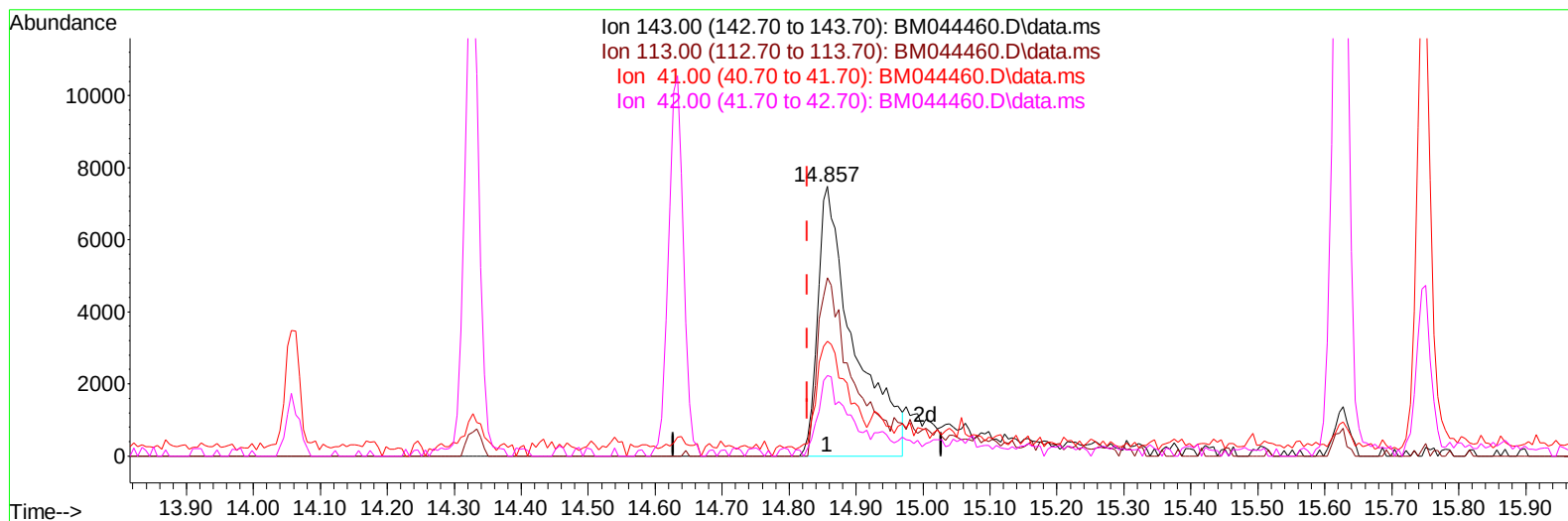
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022824\
 Data File : BM044460.D
 Acq On : 01 Mar 2024 04:15
 Operator : MA/JU
 Sample : P1501-01
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1098

Manual IntegrationsAPPROVED

Quant Time: Mar 01 04:44:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 01 01:46:17 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/01/2024
 Supervised By :mohammad ahmed 03/02/2024



TIC: BM044460.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.857min (+ 0.030) 3.25 ng/ul

response 27878

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	66.04
41.00	38.20	42.46
42.00	25.30	29.88

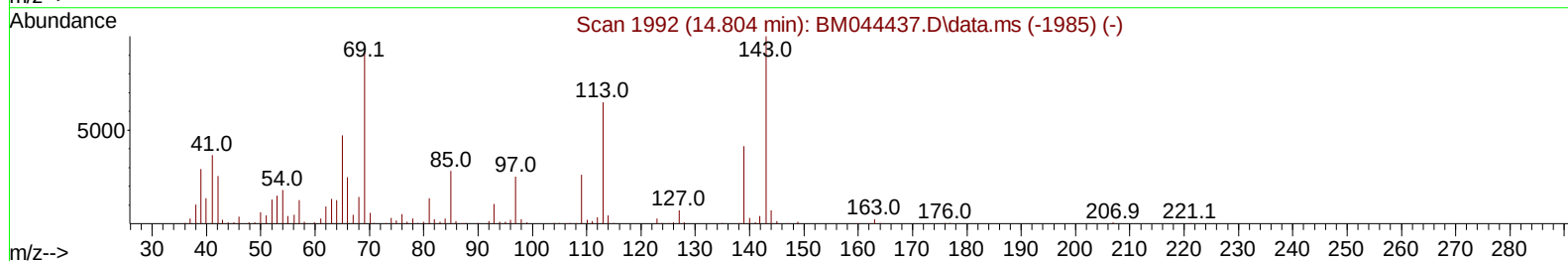
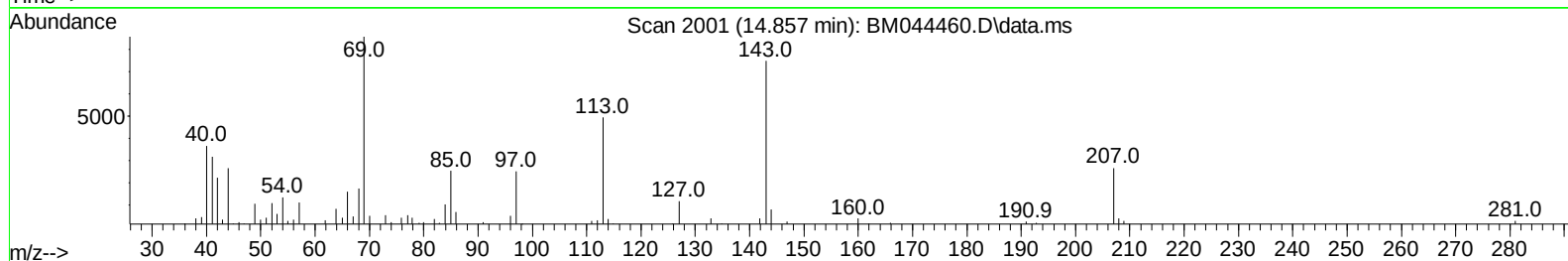
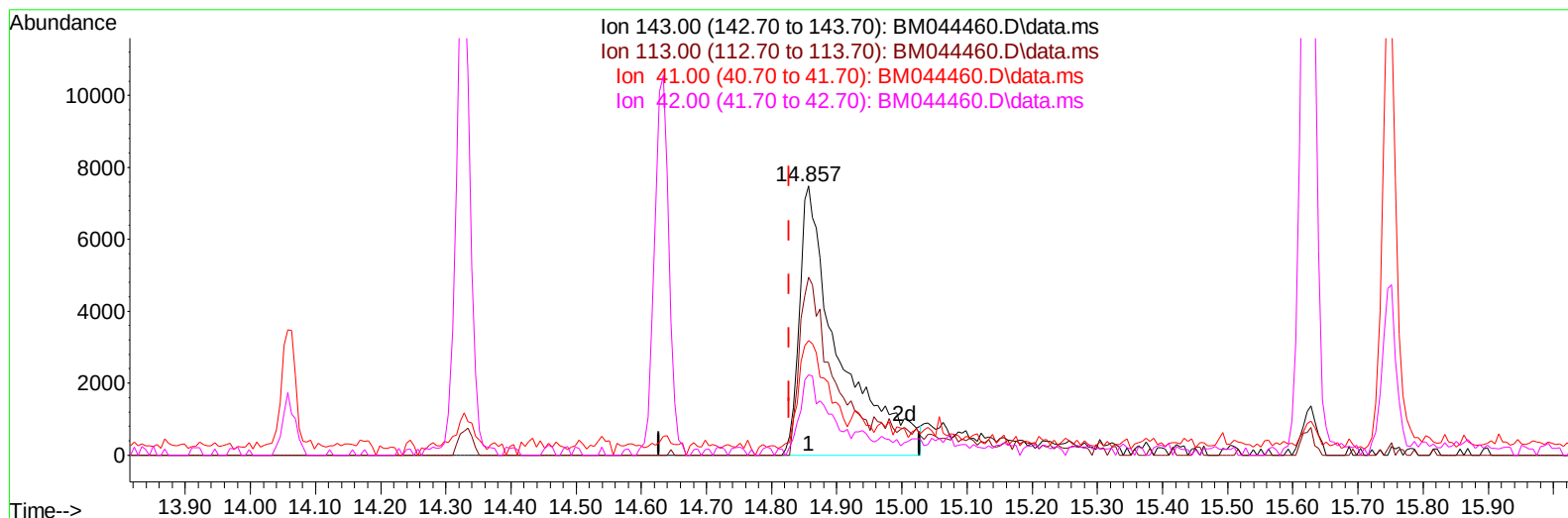
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(54) 4-Nitrophenol-d4 (S)

14.857min (+ 0.030) 3.66 ng/ul m

response 31429

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.00	66.04
41.00	38.20	42.46
42.00	25.30	29.88

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.993	152	208789	20.000	ng/ul	0.01
20) Naphthalene-d8	10.804	136	920499	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.633	164	558956	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.380	188	1099655	20.000	ng/ul	0.01
79) Chrysene-d12	21.562	240	881045	20.000	ng/ul	0.00
88) Perylene-d12	24.009	264	976377	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	25485	3.851	ng/uL	0.00
4) Pyridine-d5	3.817	84	149953	8.297	ng/ul	0.02
7) Phenol-d5	7.157	99	126668	6.333	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	423474	32.695	ng/ul	0.01
11) 2-Chlorophenol-d4	7.522	132	391055	25.959	ng/ul	0.02
15) 4-Methylphenol-d8	8.710	113	235077	15.533	ng/ul	0.01
21) Nitrobenzene-d5	9.169	128	261528	34.943	ng/ul	0.01
24) 2-Nitrophenol-d4	9.893	143	266270	34.871	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.428	165	412753	30.643	ng/ul	0.02
31) 4-Chloroaniline-d4	10.957	131	537548	23.762	ng/ul	0.02
46) Dimethylphthalate-d6	14.057	166	1478744	34.302	ng/ul	0.01
49) Acenaphthylene-d8	14.328	160	1755346	35.391	ng/ul	0.01
54) 4-Nitrophenol-d4	14.857	143	31429m	3.664	ng/ul	0.03
60) Fluorene-d10	15.622	176	1332304	36.514	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.751	200	208530	30.562	ng/ul	0.01
73) Anthracene-d10	17.480	188	2222795	42.393	ng/ul	0.01
81) Pyrene-d10	19.768	212	2439812	46.030	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.851	264	2162184	43.217	ng/ul	0.01

Target Compounds Qvalue

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

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