

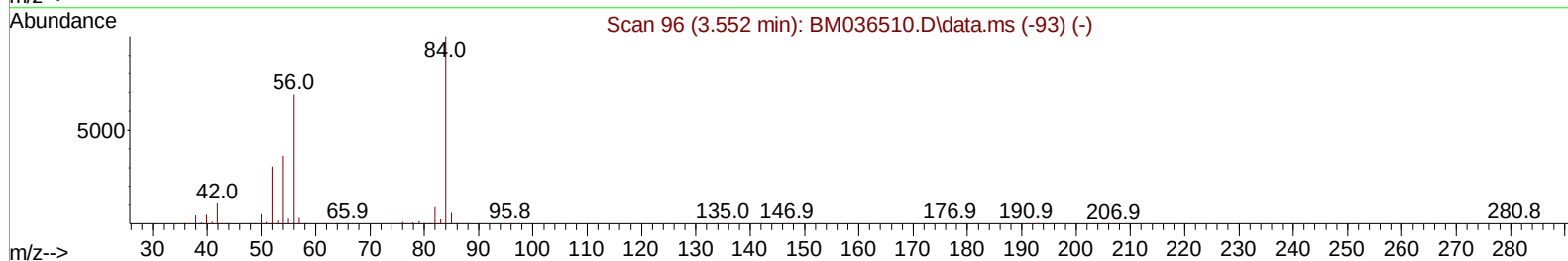
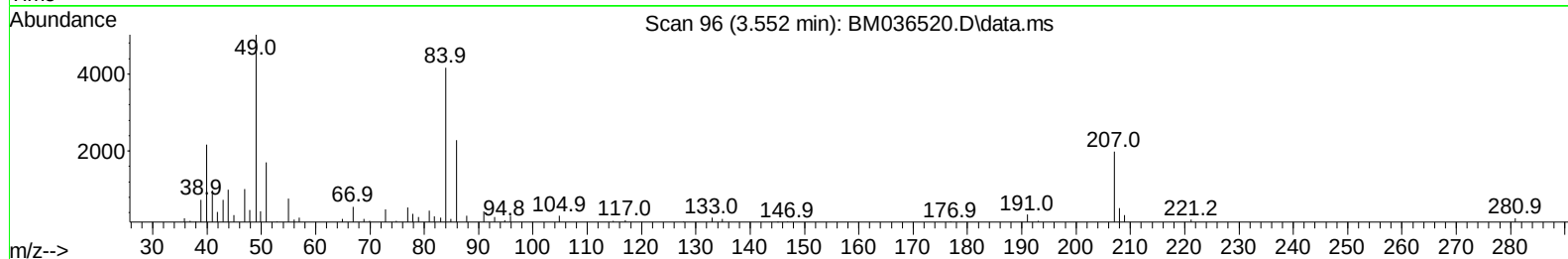
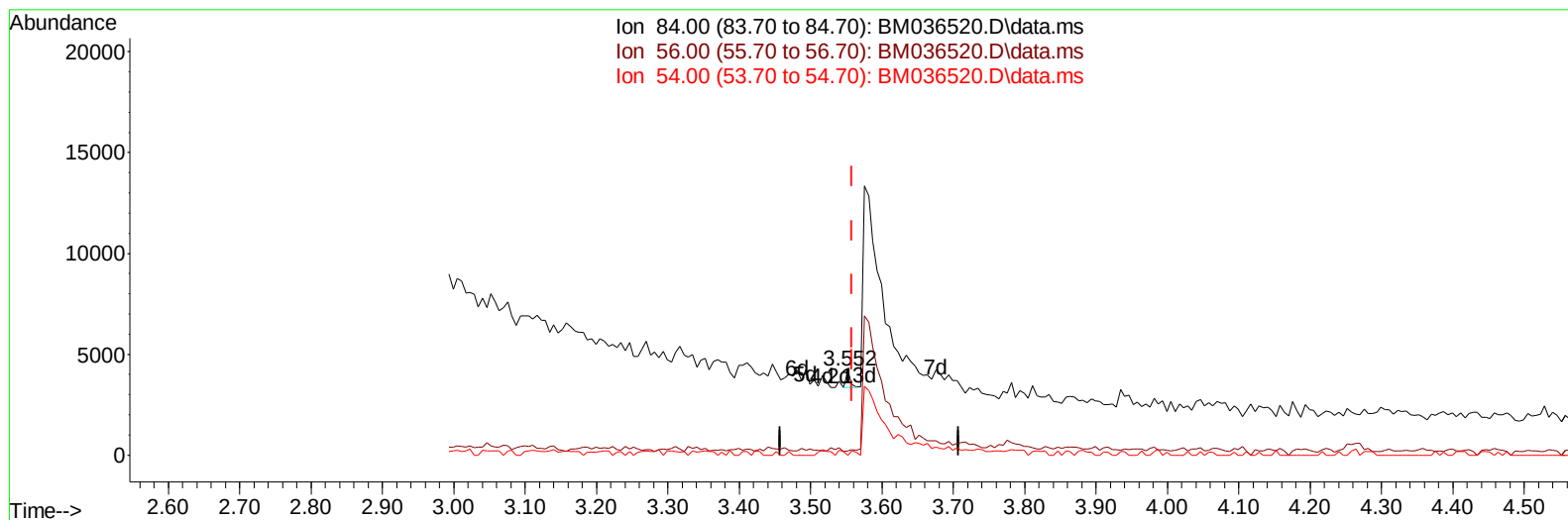
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090822\
 Data File : BM036520.D
 Acq On : 07 Sep 2022 21:49
 Operator : CG/JU
 Sample : N4483-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW353

Manual IntegrationsAPPROVED

Quant Time: Sep 08 00:55:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 02 01:33:23 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/08/2022
 Supervised By :mohammad ahmed 09/19/2022



TIC: BM036520.D\data.ms

(4) Pyridine-d5 (S)

3.552min (-0.006) 0.04 ng/ul

response 359

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.90	5.21#
54.00	32.70	0.00#
0.00	0.00	0.00

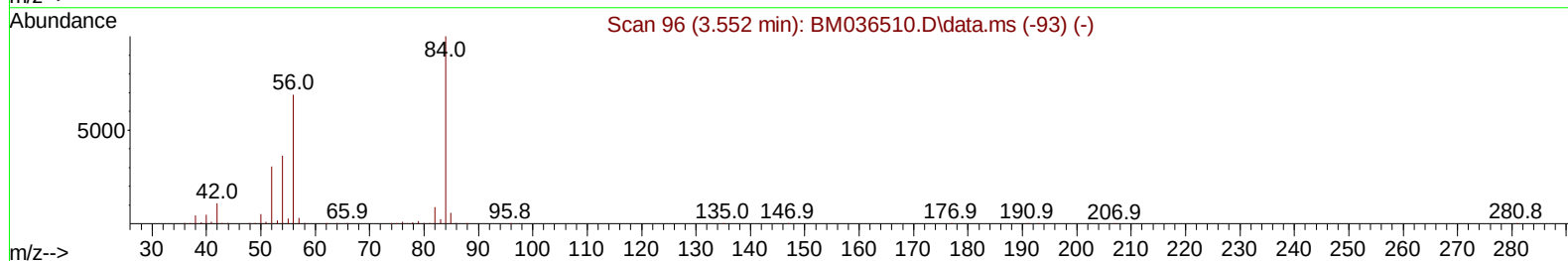
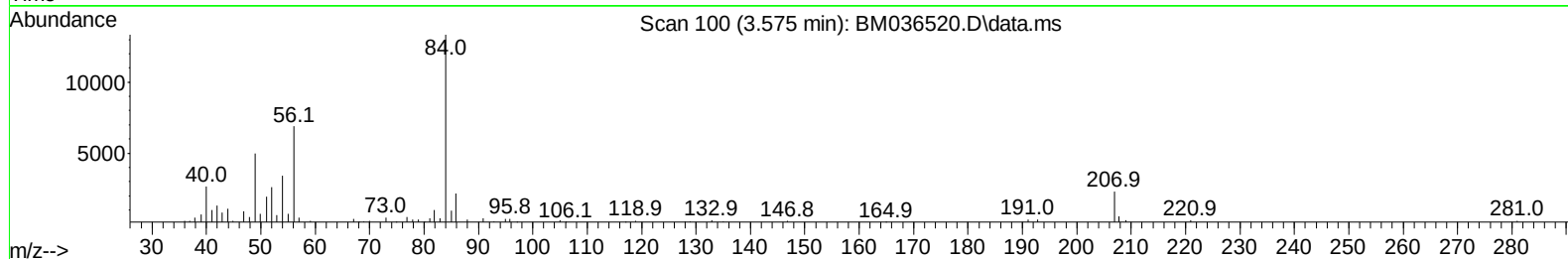
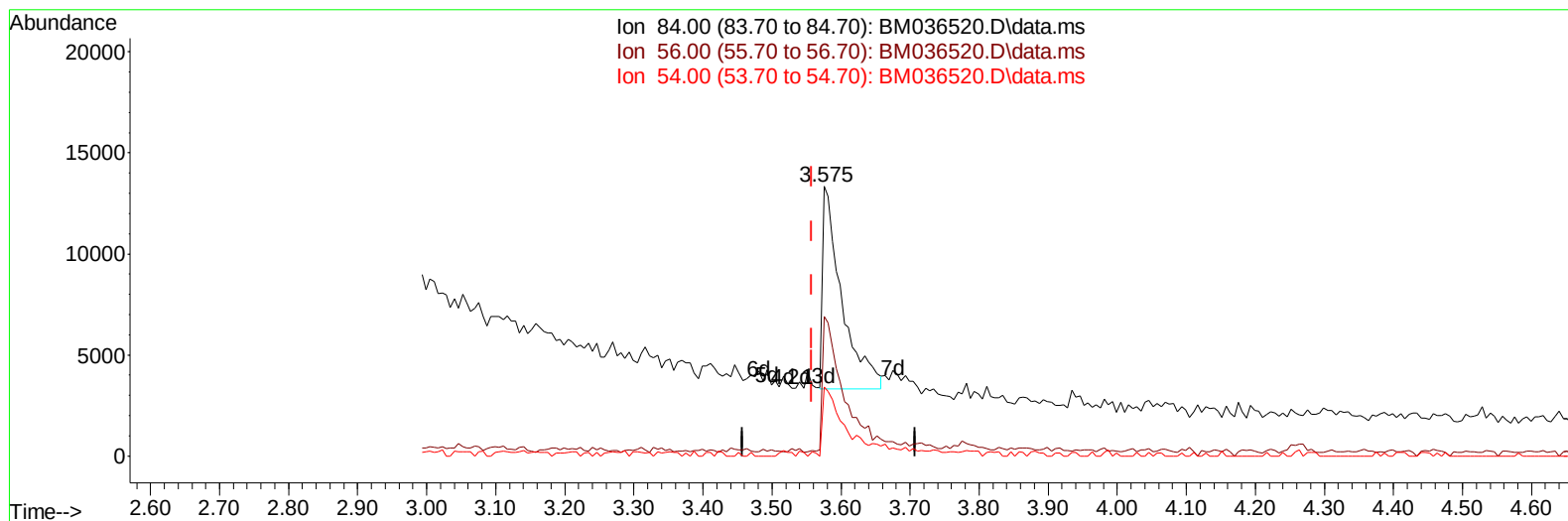
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(4) Pyridine-d5 (S)

3.575min (+ 0.018) 1.99 ng/ul m

response 19255

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.90	51.68#
54.00	32.70	25.64#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	133837	20.000	ng/ul	0.00
20) Naphthalene-d8	10.516	136	631434	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.375	164	410636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	895473	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	973023	20.000	ng/ul	0.00
88) Perylene-d12	23.598	264	934621	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	13028	3.710	ng/uL	0.00
4) Pyridine-d5	3.575	84	19255m	1.988	ng/ul	0.02
7) Phenol-d5	6.899	99	328562	28.761	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	208481	28.992	ng/ul	0.00
11) 2-Chlorophenol-d4	7.252	132	264708	29.491	ng/ul	0.00
15) 4-Methylphenol-d8	8.446	113	268481	31.475	ng/ul	0.00
21) Nitrobenzene-d5	8.893	128	141547	29.566	ng/ul	0.00
24) 2-Nitrophenol-d4	9.610	143	150659	31.025	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.146	165	267431	30.821	ng/ul	0.00
31) 4-Chloroaniline-d4	10.669	131	287291	21.358	ng/ul	0.00
46) Dimethylphthalate-d6	13.798	166	904923	34.833	ng/ul	0.00
49) Acenaphthylene-d8	14.063	160	1056538	32.608	ng/ul	0.00
54) 4-Nitrophenol-d4	14.604	143	159422	32.387	ng/ul	0.00
60) Fluorene-d10	15.369	176	840828	35.818	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	109319	23.746	ng/ul	0.00
73) Anthracene-d10	17.216	188	1422115	35.716	ng/ul	0.00
81) Pyrene-d10	19.516	212	1824633	36.150	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.451	264	1780362	38.852	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	6.934	94	14710	1.238	ng/ul#	88
16) Acetophenone	8.557	105	18886	1.377	ng/ul	97
80) Fluoranthene	19.180	202	61322	1.005	ng/ul	99

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

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