

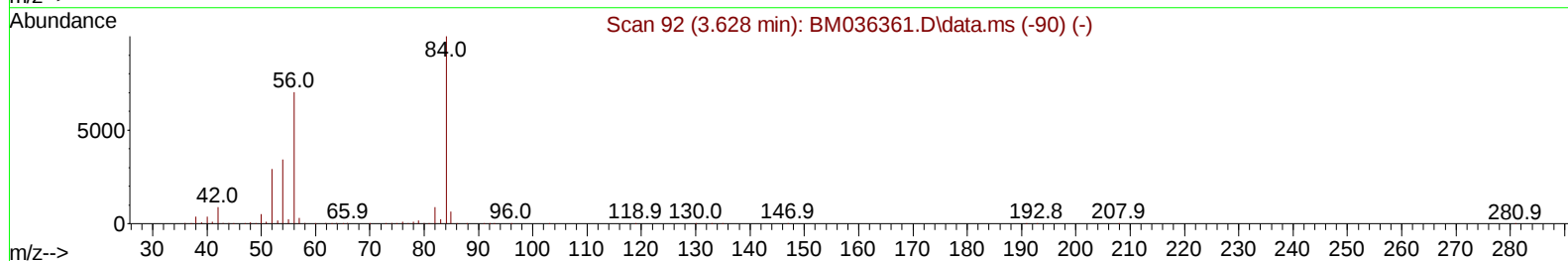
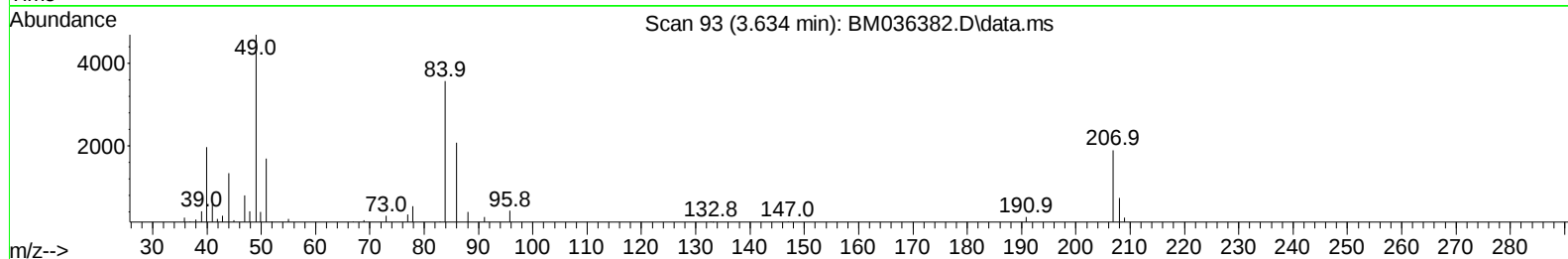
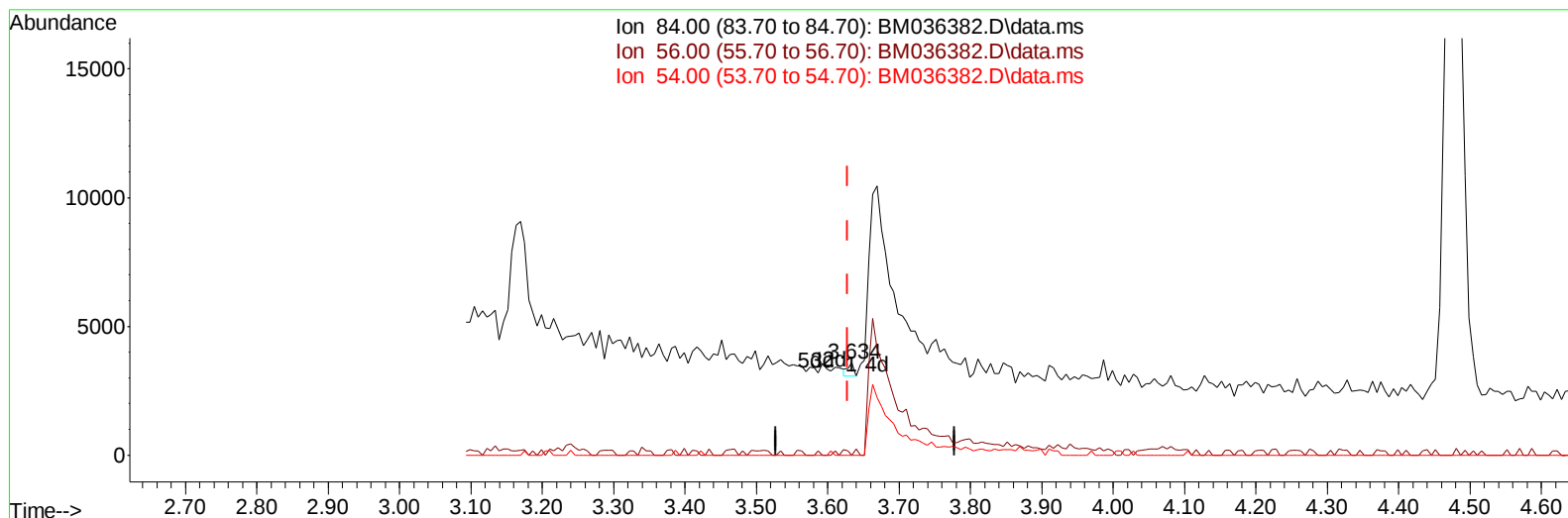
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082222\
 Data File : BM036382.D
 Acq On : 23 Aug 2022 09:11
 Operator : CG/JU
 Sample : N4231-06DL 10X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0KN4DL

Manual IntegrationsAPPROVED

Quant Time: Aug 24 00:55:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 23 00:55:01 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/24/2022
 Supervised By :Sohil Jodhani 08/24/2022



TIC: BM036382.D\data.ms

(4) Pyridine-d5 (S)

3.634min (+ 0.006) 0.02 ng/u1

response 277

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

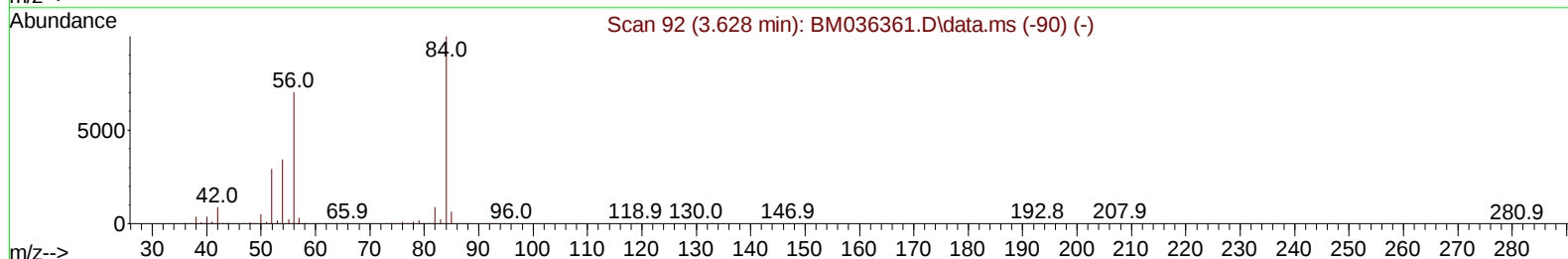
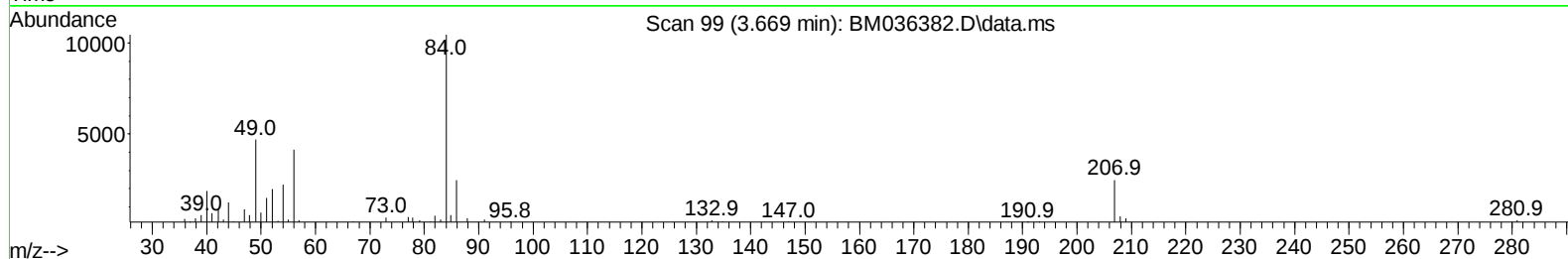
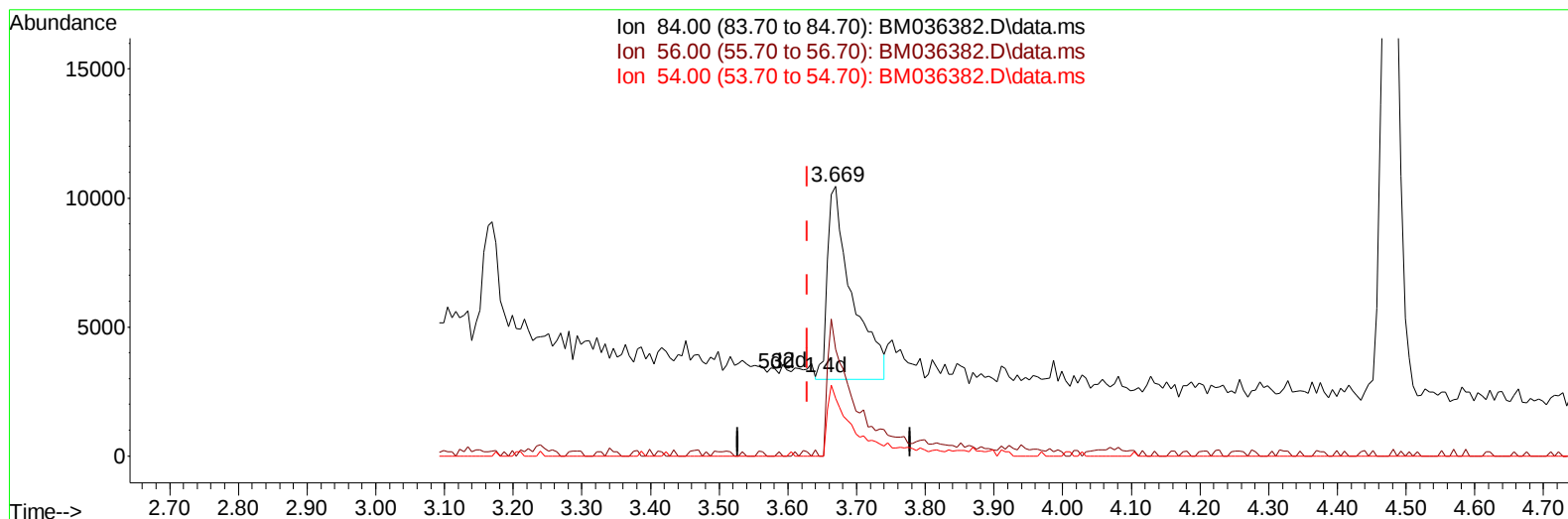
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(4) Pyridine-d5 (S)

3.669min (+ 0.041) 1.16 ng/ul m

response 18614

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.90	39.56#
54.00	32.70	21.36#
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	7.787	152	222107	20.000	ng/ul	0.00
20) Naphthalene-d8	10.586	136	1012352	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.433	164	645472	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.174	188	1210077	20.000	ng/ul	0.00
79) Chrysene-d12	21.362	240	825046	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	774438	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	2160	0.371	ng/uL	0.00
4) Pyridine-d5	3.669	84	18614m	1.158	ng/ul	0.04
7) Phenol-d5	6.981	99	15368	0.811	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	57782	4.842	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	46847	3.145	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	28852	2.038	ng/ul	0.01
21) Nitrobenzene-d5	8.963	128	28613	3.728	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	25416	3.264	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.239	165	43669	3.139	ng/ul	0.02
31) 4-Chloroaniline-d4	10.757	131	69727	3.233	ng/ul	0.02
46) Dimethylphthalate-d6	13.857	166	197687	4.841	ng/ul	0.00
49) Acenaphthylene-d8	14.127	160	212657	4.175	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.427	176	178959	4.850	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.280	188	248993	4.628	ng/ul	0.00
81) Pyrene-d10	19.574	212	239015	5.585	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	178443	4.700	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	75972	12.377	ng/uL	94

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

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