

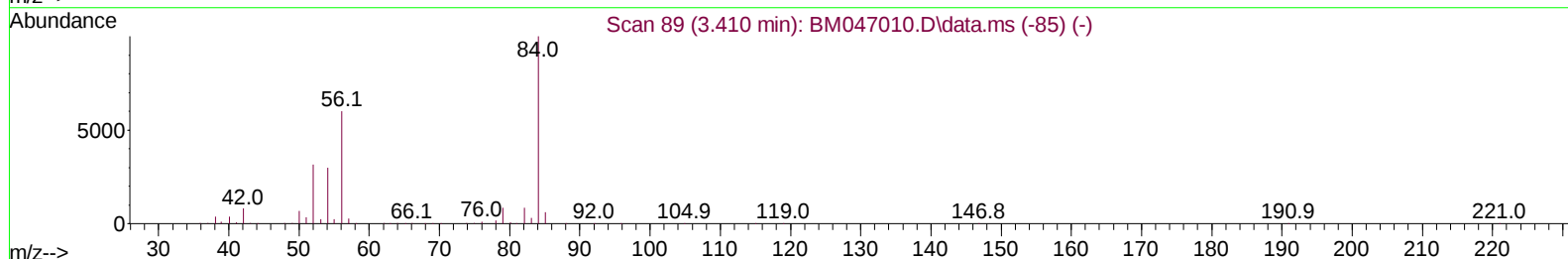
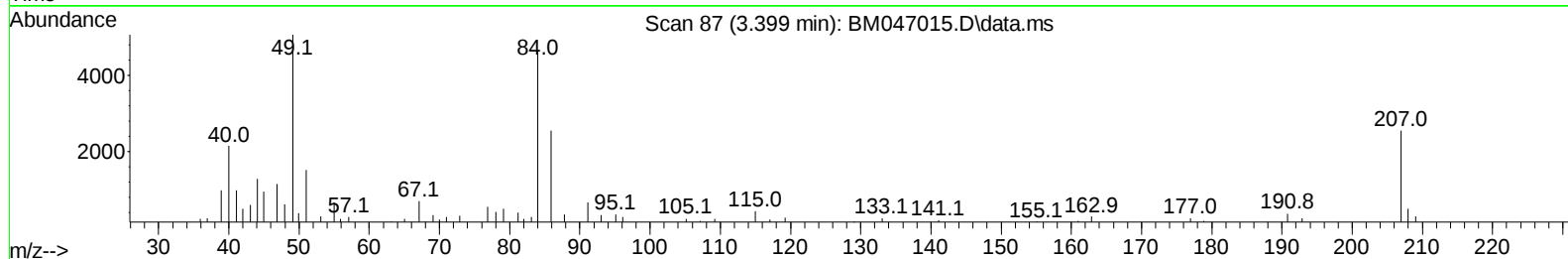
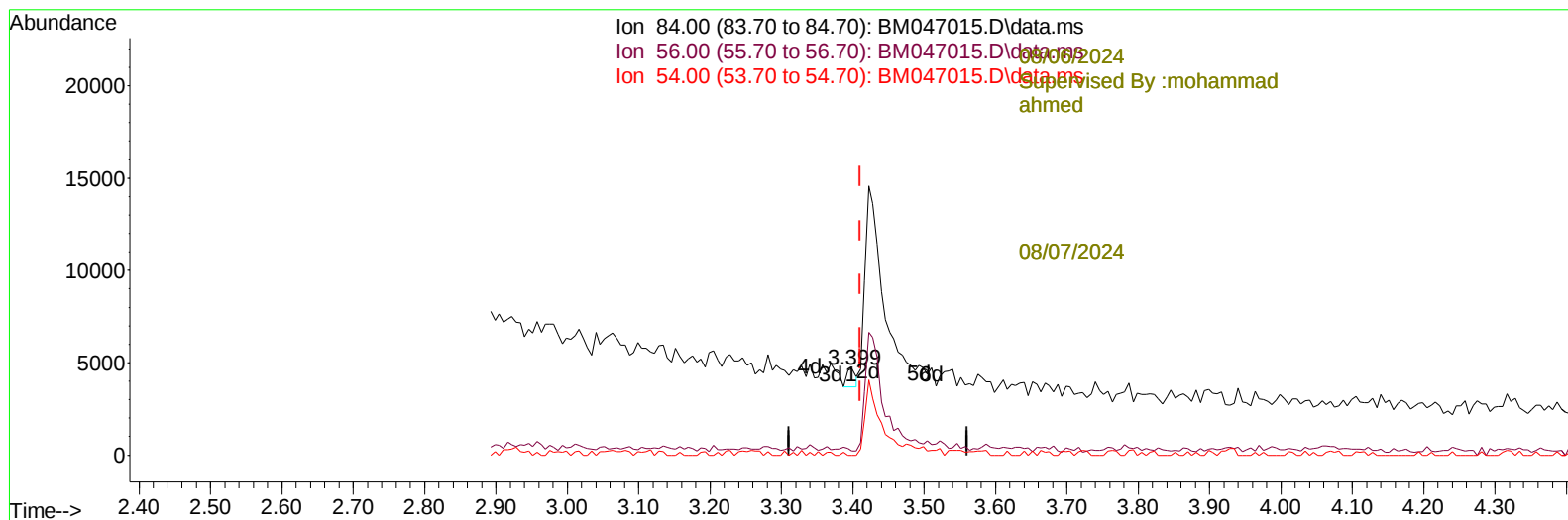
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047015.D
Acq On : 06 Aug 2024 07:32
Operator : RC/JU
Sample : P3382-02
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E20R8

Manual IntegrationsAPPROVED

Quant Time: Aug 06 08:13:18 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 06 04:53:02 2024
Response via : Initial Calibration

Reviewed By :Jagrut
Upadhyay



TIC: BM047015.D\data.ms

(4) Pyridine-d5 (S)

3.399min (-0.012) 0.04 ng/u1

response 845

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.50	4.96#
54.00	34.00	0.00#
0.00	0.00	0.00

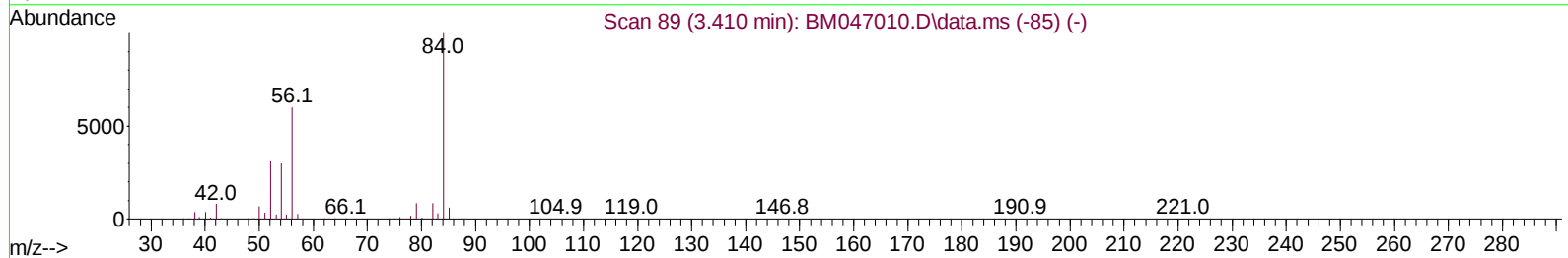
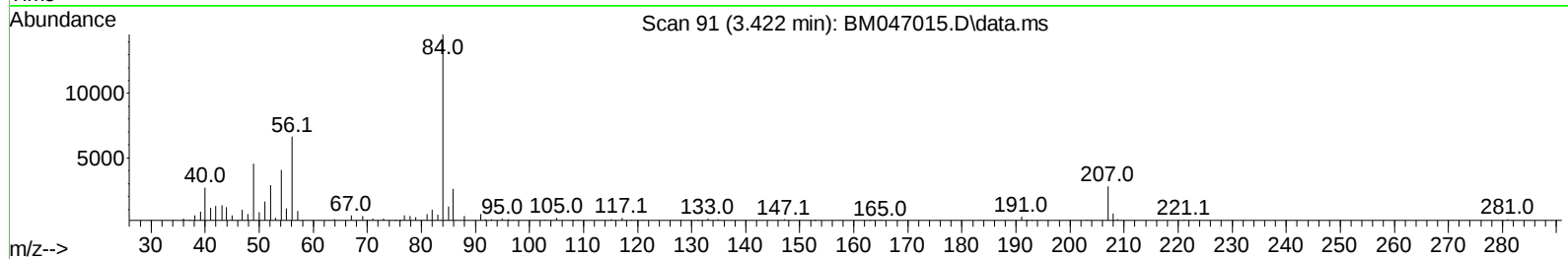
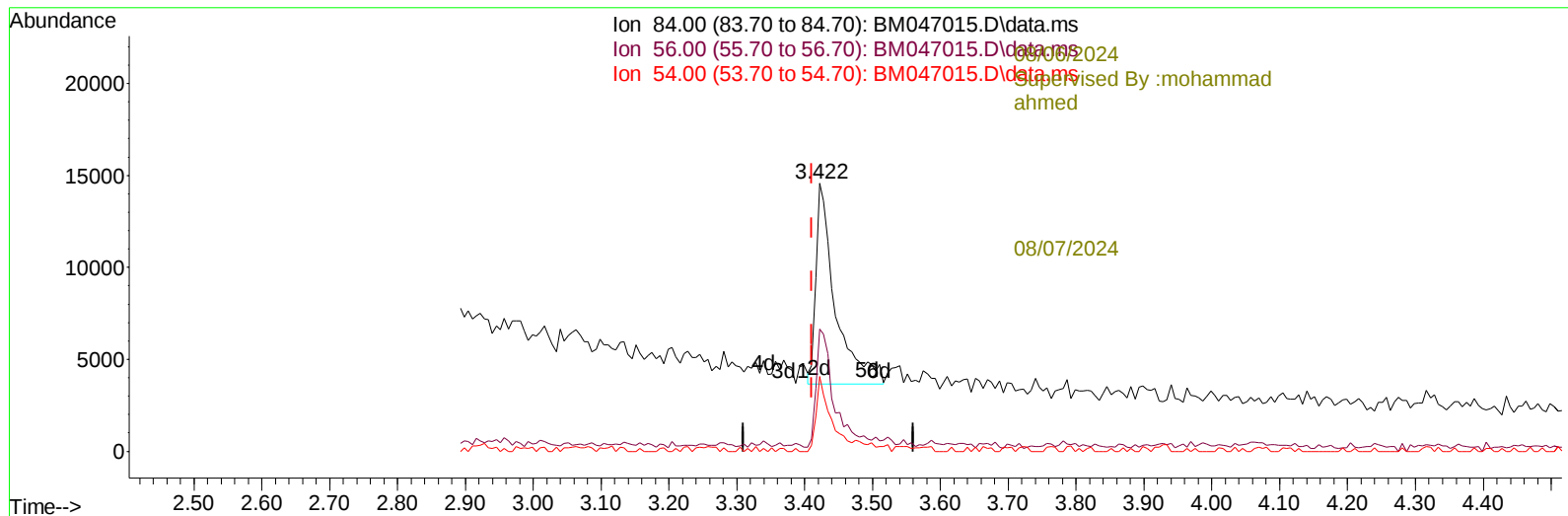
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(4) Pyridine-d5 (S)

3.422min (+ 0.012) 1.07 ng/ul m

response 21559

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.50	45.63#
54.00	34.00	27.85
0.00	0.00	0.00

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 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
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 E20R8

Manual IntegrationsAPPROVED

Quant Time: Aug 06 08:13:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 06 04:53:02 2024
 Response via : Initial Calibration

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----08/06/2024							
Supervised By :mohammad ali med							
-----08/07/2024							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.392	152		293805	20.000	ng/ul	0.00
20) Naphthalene-d8	10.151	136		1181445	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.051	164		803638	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.809	188		1696252	20.000	ng/ul	0.00
79) Chrysene-d12	21.044	240		1466255	20.000	ng/ul	0.00
88) Perylene-d12	23.756	264		1641308	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.022	96		10525	1.419	ng/uL	0.00
4) Pyridine-d5	3.422	84		21559m	1.066	ng/ul	0.01
7) Phenol-d5	6.598	99		243641	10.916	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67		153380	10.444	ng/ul	0.00
11) 2-Chlorophenol-d4	6.934	132		199981	10.578	ng/ul	0.00
15) 4-Methylphenol-d8	8.116	113		180914	10.232	ng/ul	0.00
21) Nitrobenzene-d5	8.539	128		95037	10.025	ng/ul	0.00
24) 2-Nitrophenol-d4	9.251	143		108764	10.505	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.786	165		213477	10.740	ng/ul	0.00
31) 4-Chloroaniline-d4	10.304	131		241157	9.231	ng/ul	0.00
46) Dimethylphthalate-d6	13.480	166		645034	10.497	ng/ul	0.00
49) Acenaphthylene-d8	13.739	160		740479	10.796	ng/ul	0.00
54) 4-Nitrophenol-d4	14.292	143		100461	8.456	ng/ul	0.00
60) Fluorene-d10	15.057	176		580068	11.035	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.192	200		71204	6.644	ng/ul	0.00
73) Anthracene-d10	16.909	188		854062	10.621	ng/ul	0.00
81) Pyrene-d10	19.221	212		843594	10.051	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264		608966	7.046	ng/ul	0.00
Target Compounds							
						Qvalue	
16) Acetophenone	8.210	105		74409	2.674	ng/ul	98
47) Dimethylphthalate	13.527	163		183454	2.975	ng/ul	99
80) Fluoranthene	18.886	202		192253	1.917	ng/ul	99
82) Pyrene	19.250	202		118716	1.111	ng/ul	96
90) Benzo(b)fluoranthene	22.909	252		131763	1.298	ng/ul#	88

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

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