

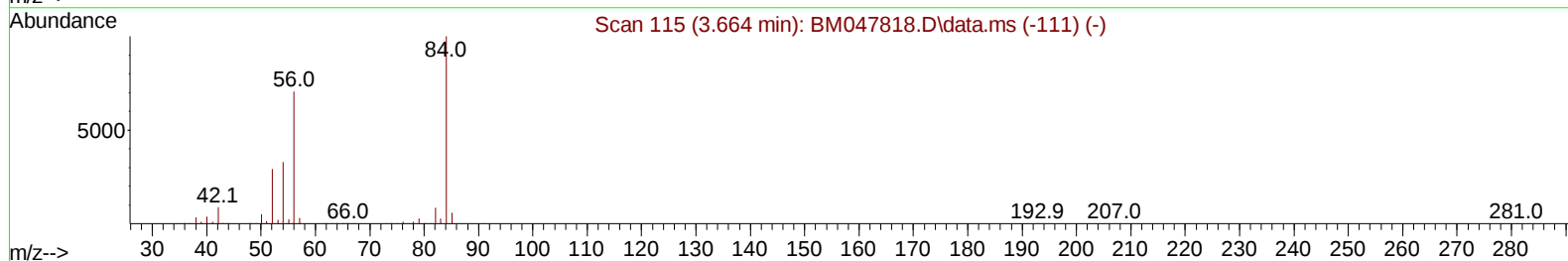
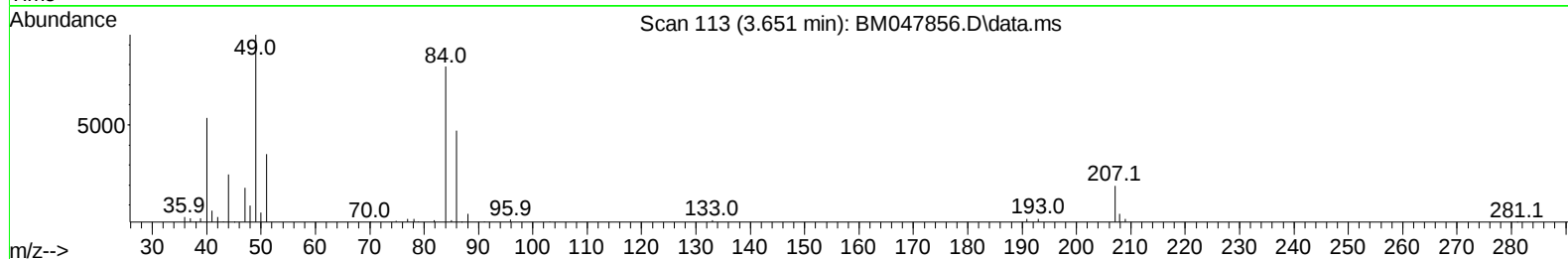
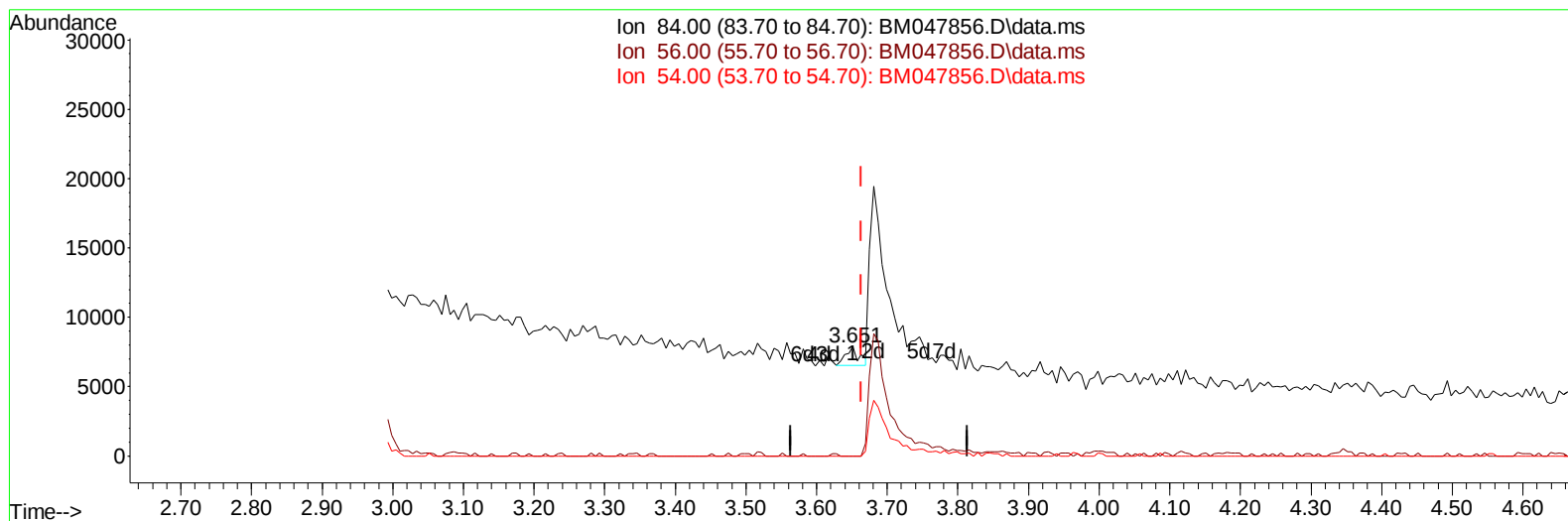
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100324\
 Data File : BM047856.D
 Acq On : 05 Oct 2024 00:13
 Operator : RC/JU
 Sample : P4017-03MEDL 20X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDC42MEDL

Manual IntegrationsAPPROVED

Quant Time: Oct 05 00:55:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 08:32:31 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/05/2024
 Supervised By :mohammad ahmed 10/06/2024



TIC: BM047856.D\data.ms

(4) Pyridine-d5 (S)

3.651min (-0.012) 0.09 ng/u1

response 1706

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	75.30	0.00#
54.00	36.10	0.00#
0.00	0.00	0.00

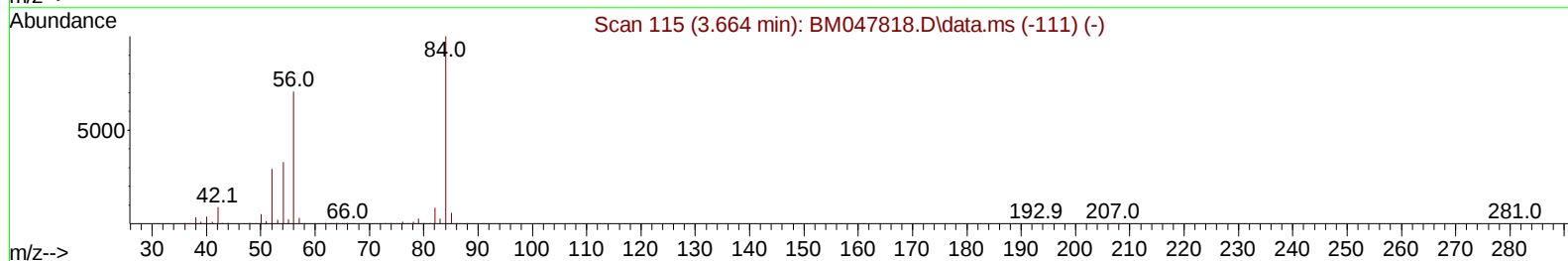
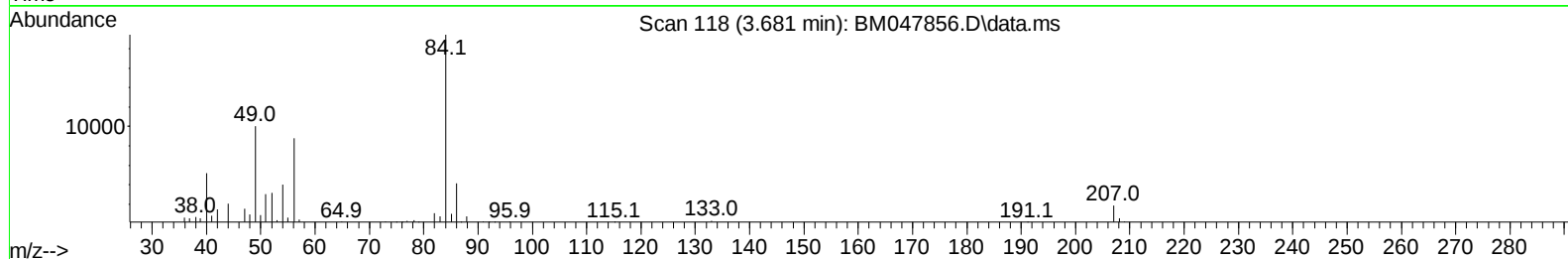
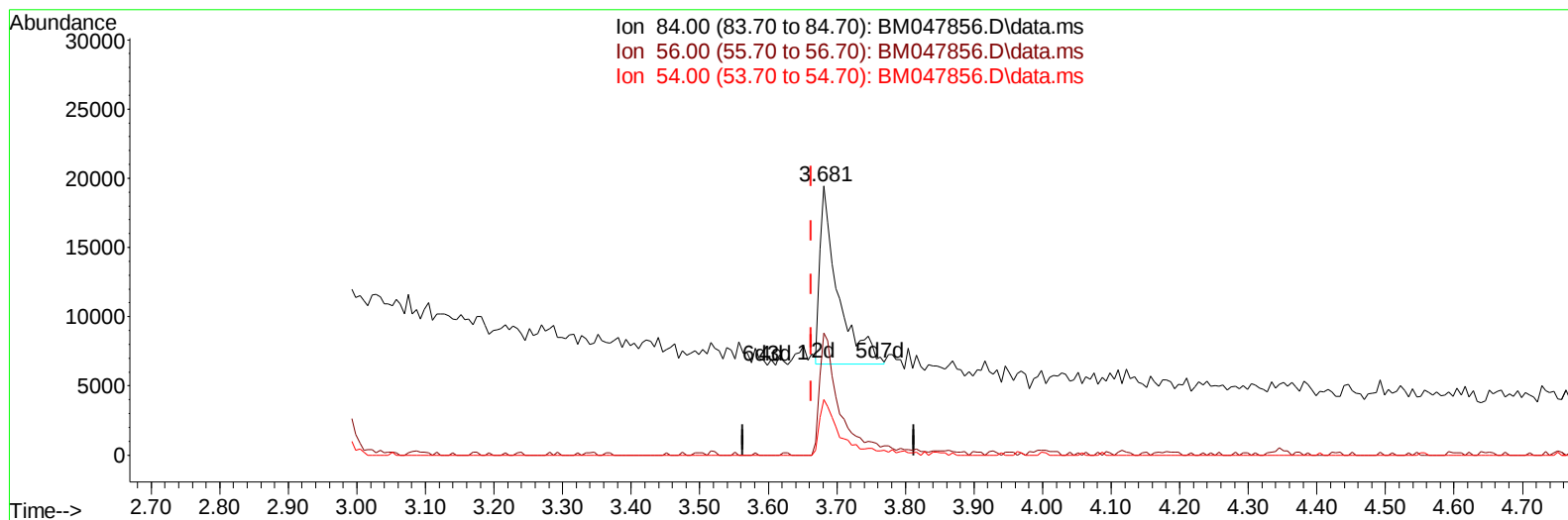
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(4) Pyridine-d5 (S)

3.681min (+ 0.018) 1.27 ng/u1 m

response 23491

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	75.30	45.30#
54.00	36.10	20.63#
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/05/2024
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Quant Time: Oct 05 00:57:15 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	204565	20.000	ng/ul	0.00
20) Naphthalene-d8	10.592	136	815224	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.433	164	492914	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.168	188	1037315	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240	956697	20.000	ng/ul	0.00
88) Perylene-d12	24.391	264	1081131	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.681	84	23491m	1.267	ng/ul	0.02
7) Phenol-d5	6.969	99	25190	1.302	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	19258	1.403	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	22066	1.568	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	19359	1.346	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	9637	1.485	ng/ul	0.00
24) 2-Nitrophenol-d4	9.675	143	8368	1.345	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	17775	1.431	ng/ul	0.00
31) 4-Chloroaniline-d4	10.727	131	27917	1.453	ng/ul	0.00
46) Dimethylphthalate-d6	13.839	166	67367	1.885	ng/ul	0.00
49) Acenaphthylene-d8	14.127	160	76319	1.801	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.421	176	59993	1.938	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	3605	0.594	ng/ul	0.00
73) Anthracene-d10	17.268	188	90638	1.777	ng/ul	0.00
81) Pyrene-d10	19.556	212	104179	1.919	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.185	264	92480	1.649	ng/ul	0.00
Target Compounds						
						Qvalue
58) 2,3,4,6-Tetrachlorophenol	15.057	232	33178	4.146	ng/ul	98
71) Pentachlorophenol	16.827	266	454086	57.390	ng/ul	100

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

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