

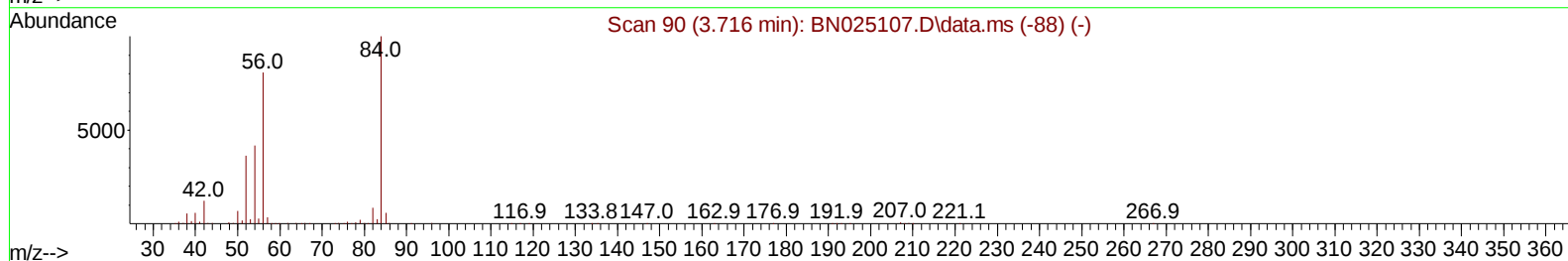
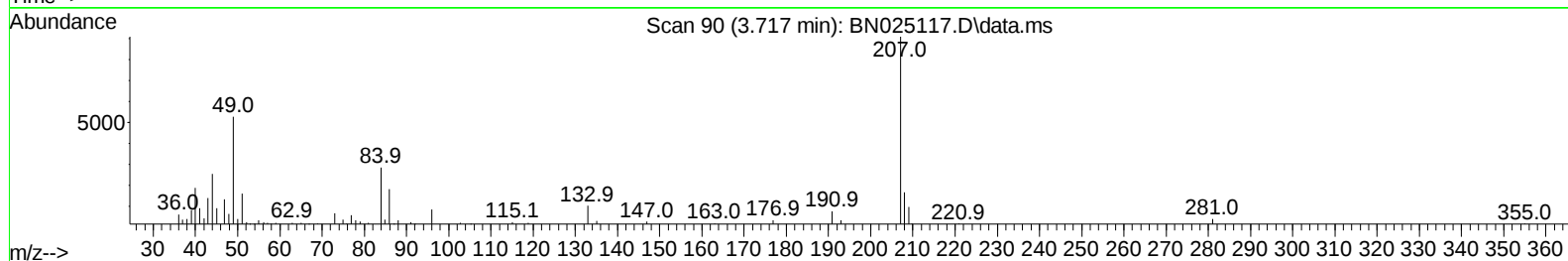
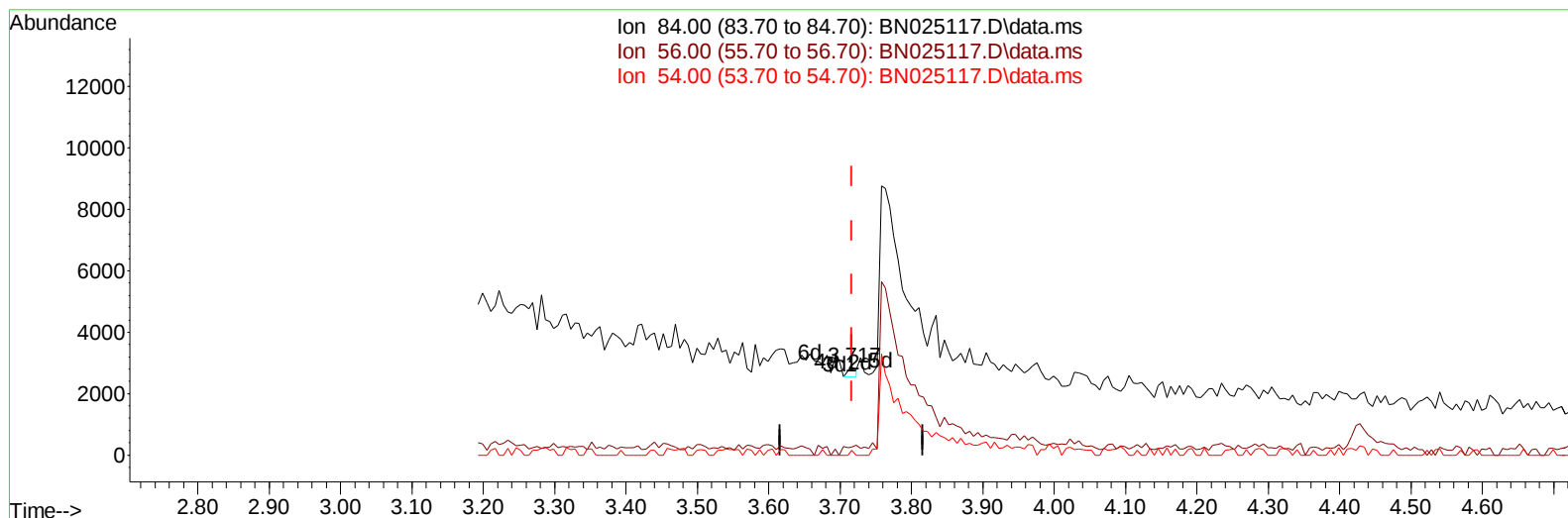
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042423\
Data File : BN025117.D
Acq On : 24 Apr 2023 16:43
Operator : CG/JU
Sample : 02440-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
COMQ6

Manual IntegrationsAPPROVED

Quant Time: Apr 25 01:26:24 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Apr 23 01:54:00 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/25/2023
Supervised By :Jagrut Upadhyay 04/27/2023



TIC: BN025117.D\data.ms

(4) Pyridine-d5 (S)

3.717min (-0.000) 0.01 ng/u1

response 256

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	9.28#
54.00	40.20	5.62#
0.00	0.00	0.00

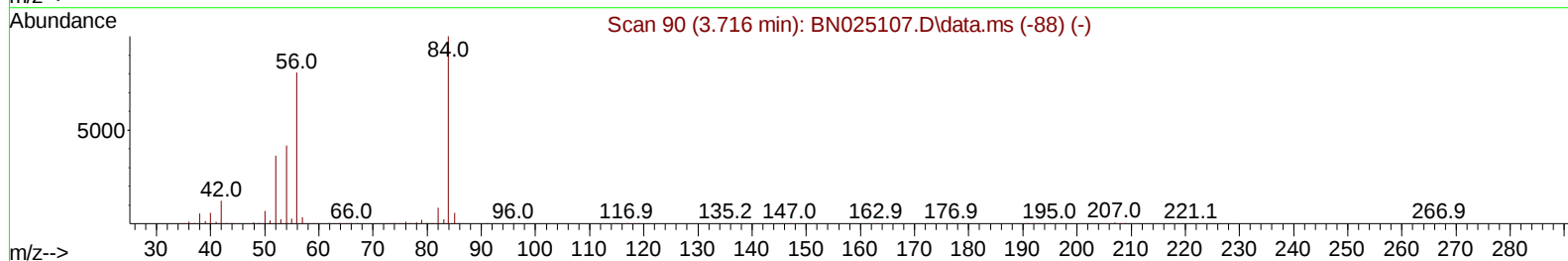
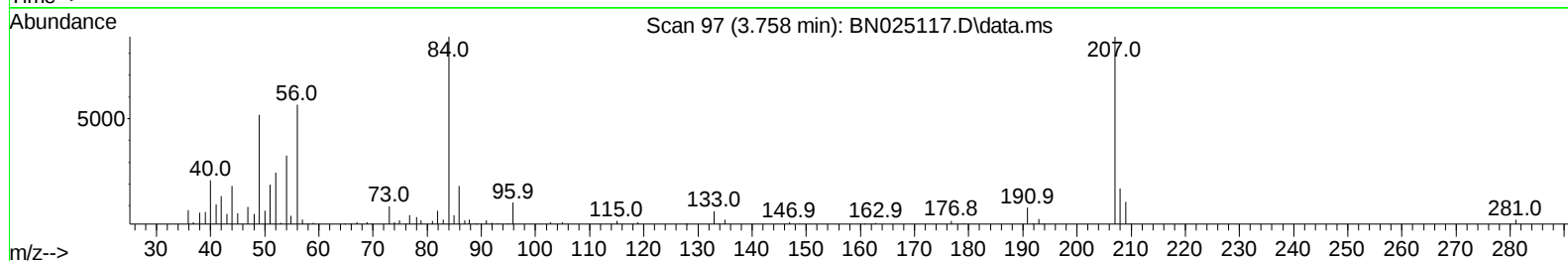
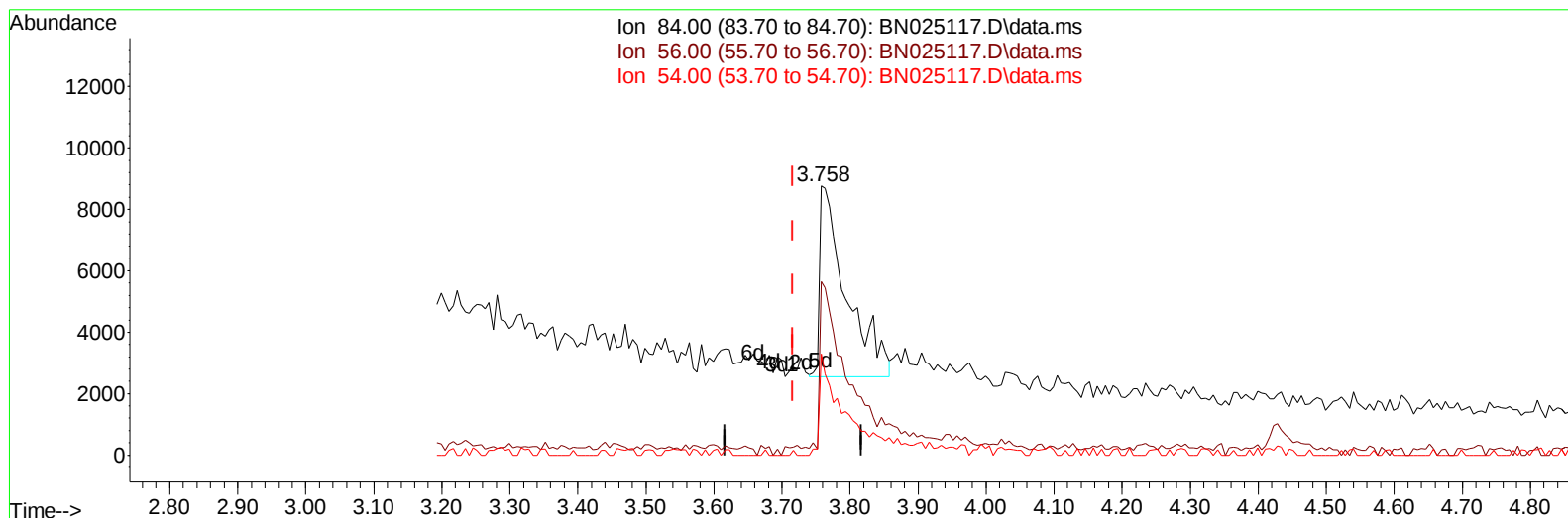
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(4) Pyridine-d5 (S)

3.758min (+ 0.041) 0.92 ng/u1 m

response 16965

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	64.39#
54.00	40.20	37.72
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	7.916	152	234581	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	1082466	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	705231	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1607166	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	1615253	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	1756096	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25688	4.168	ng/uL	0.00
4) Pyridine-d5	3.758	84	16965m	0.919	ng/ul	0.04
7) Phenol-d5	7.075	99	139049	6.080	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.240	67	428551	28.016	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	373040	22.195	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	258764	14.252	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	234738	26.783	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	253720	26.726	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	444394	26.189	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	44325	1.721	ng/ul	0.00
46) Dimethylphthalate-d6	13.969	166	1654046	30.255	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1737011	27.726	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	53653	4.788	ng/ul	0.00
60) Fluorene-d10	15.557	176	1419443	30.751	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	216834	20.731	ng/ul	0.00
73) Anthracene-d10	17.410	188	2346936	31.092	ng/ul	0.00
81) Pyrene-d10	19.704	212	3007653	32.066	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	2982783	32.485	ng/ul	0.01

Target Compounds Qvalue

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

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