

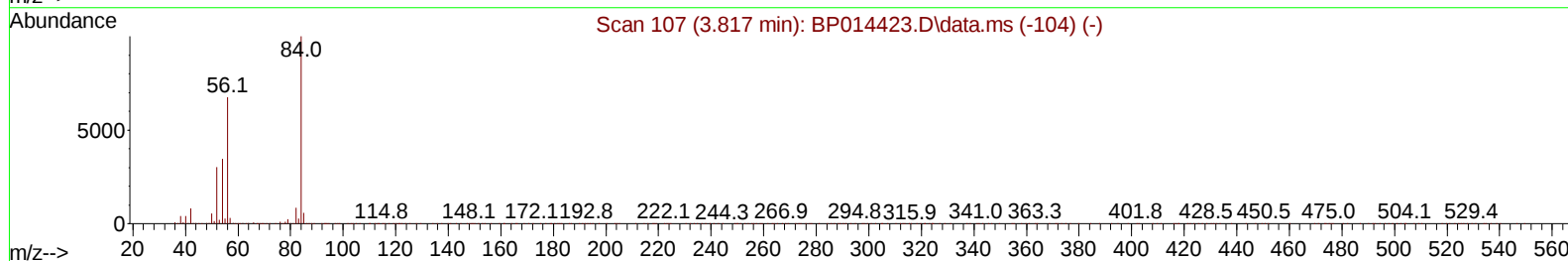
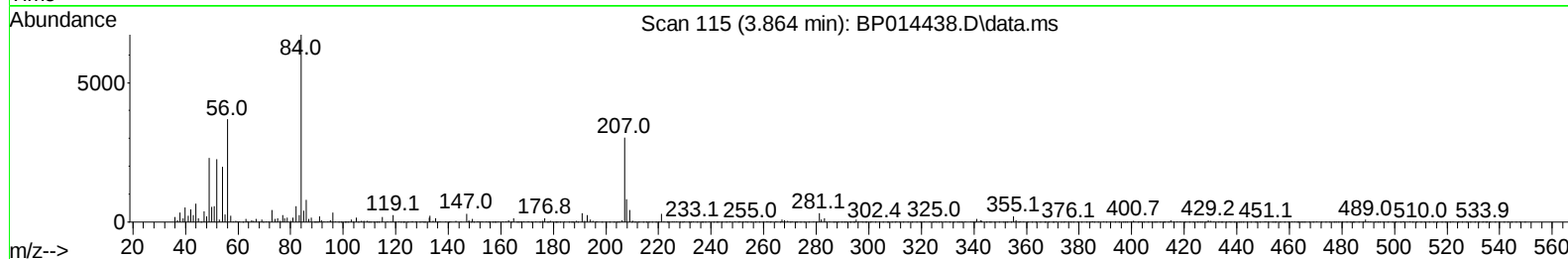
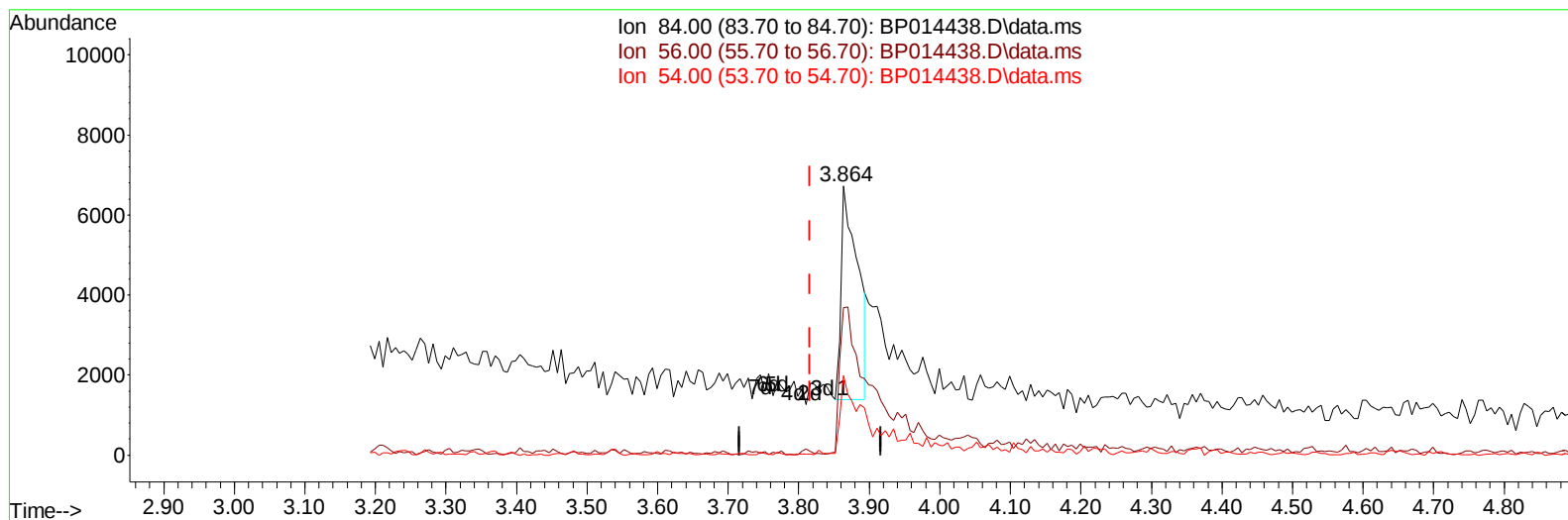
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0QG7DL

Manual IntegrationsAPPROVED

Quant Time: Mar 31 22:52:39 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 31 04:41:29 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/03/2023
Supervised By :Jagrut Upadhyay 04/03/2023



TIC: BP014438.D\data.ms

(4) Pyridine-d5 (S)

3.864min (+ 0.047) 0.79 ng/u1

response 8718

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	54.89
54.00	32.60	29.59
0.00	0.00	0.00

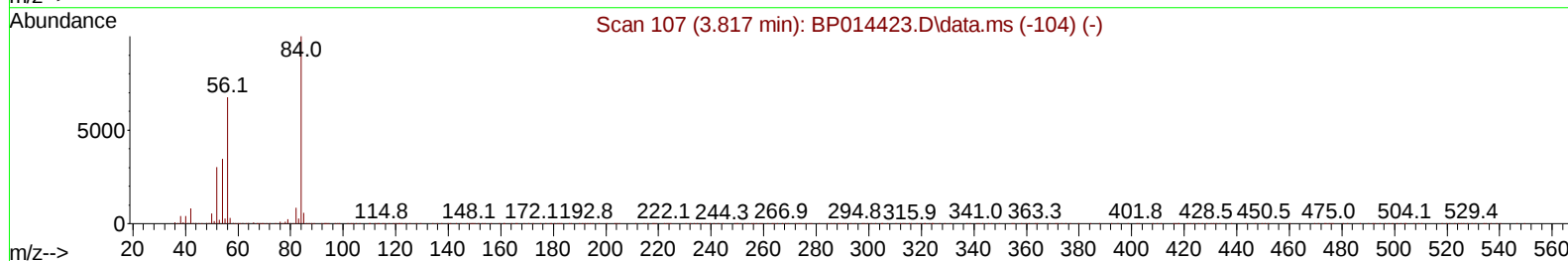
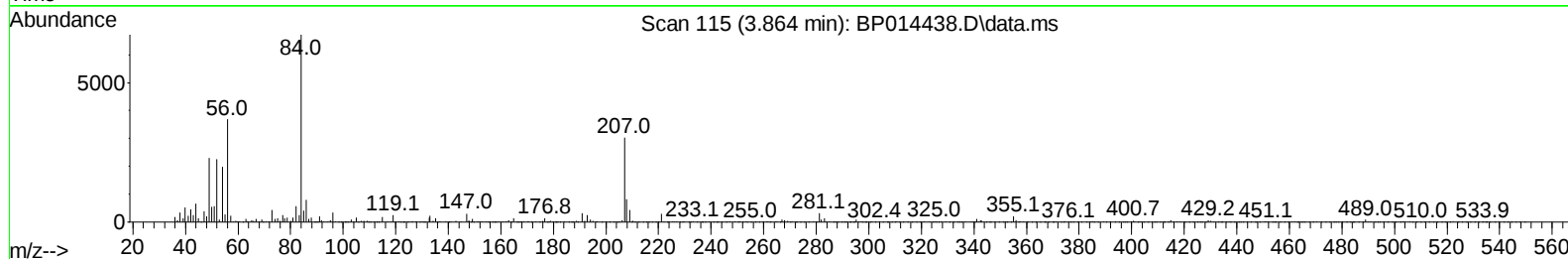
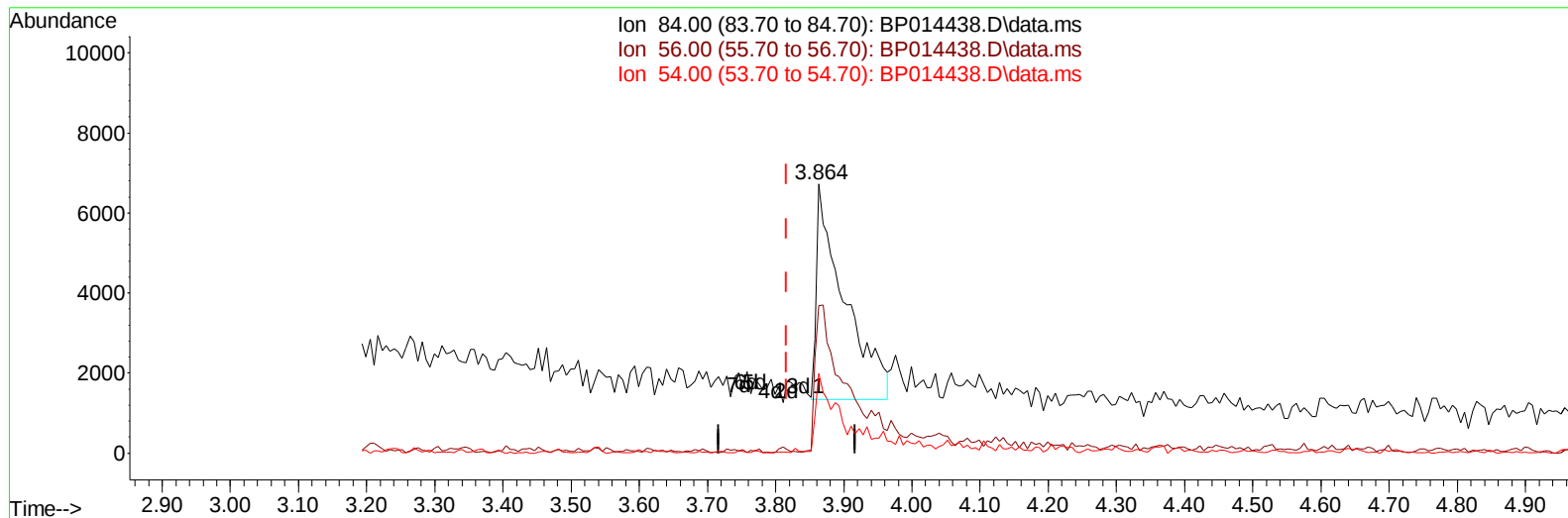
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TIC: BP014438.D\data.ms

(4) Pyridine-d5 (S)

3.864min (+ 0.047) 1.38 ng/ul m

response 15164

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	54.89
54.00	32.60	29.59
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	8.010	152	166380	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	753442	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.663	164	518313	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1118260	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	778118	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	744121	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	4358	1.016	ng/uL	0.00
4) Pyridine-d5	3.864	84	15164m	1.379	ng/ul	0.05
7) Phenol-d5	7.187	99	17558	1.144	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.340	67	69440	7.053	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	54997	4.959	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	36182	2.923	ng/ul	0.00
21) Nitrobenzene-d5	9.187	128	35758	5.879	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	41006	5.888	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	66023	5.205	ng/ul	0.01
31) 4-Chloroaniline-d4	10.993	131	51684	3.093	ng/ul	0.01
46) Dimethylphthalate-d6	14.075	166	324005	7.710	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	329293	7.062	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.657	176	268732	7.668	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	41159	5.002	ng/ul	0.00
73) Anthracene-d10	17.528	188	411662	7.607	ng/ul	0.00
81) Pyrene-d10	19.757	212	437737	8.375	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	296119	7.490	ng/ul	0.00
Target Compounds						
					Qvalue	
30) Naphthalene	10.887	128	1954702	46.576	ng/ul	99
36) 2-Methylnaphthalene	12.493	142	215341	7.571	ng/ul	98
37) 1-Methylnaphthalene	12.710	142	181531	6.436	ng/ul	94
43) 1,1'-Biphenyl	13.498	154	61582	1.471	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.298	232	158444	14.481	ng/ul	96
71) Pentachlorophenol	17.086	266	2739228	295.820	ng/ul	99
72) Phenanthrene	17.469	178	95252	1.531	ng/ul	98

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

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