

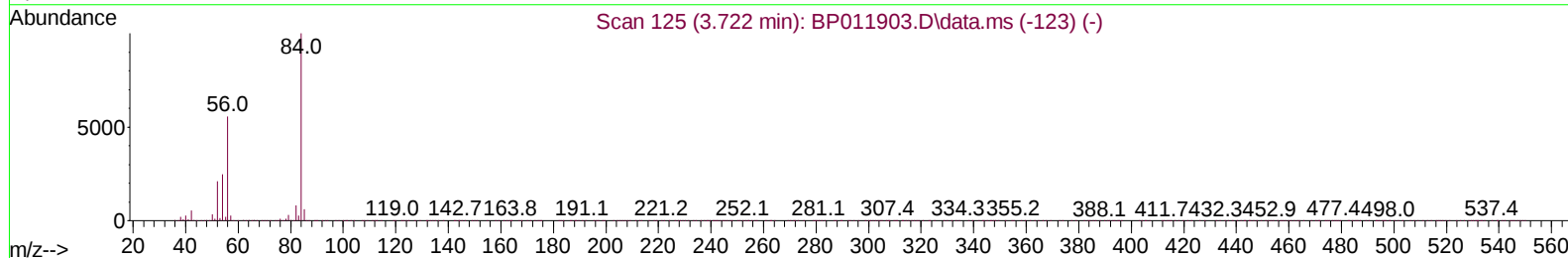
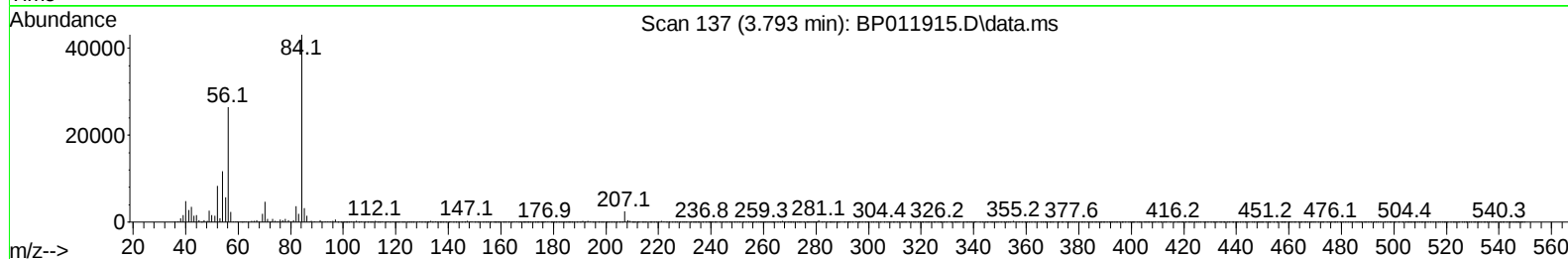
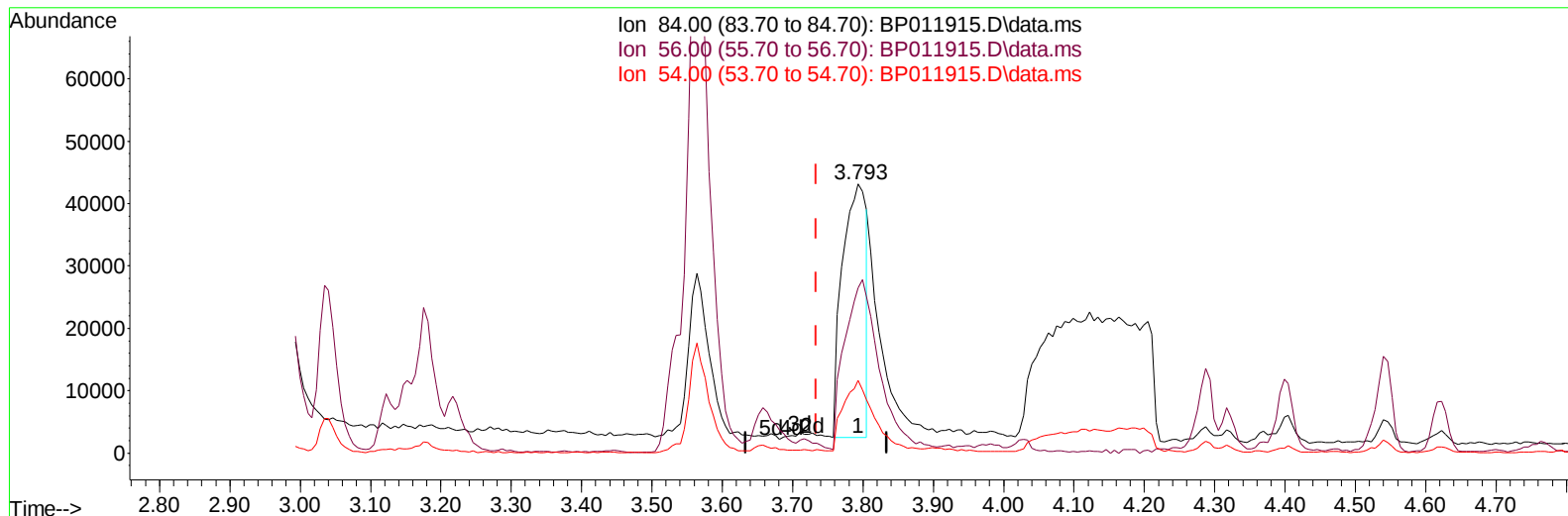
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
Data File : BP011915.D
Acq On : 04 Oct 2022 16:43
Operator : CG/JU
Sample : N4873-17
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXZ62

Manual IntegrationsAPPROVED

Quant Time: Oct 05 05:40:12 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Oct 01 00:53:59 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/05/2022
Supervised By :mohammad ahmed 10/06/2022



TIC: BP011915.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.059) 6.37 ng/u1

response 95059

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	56.10	61.35
54.00	24.80	26.96
0.00	0.00	0.00

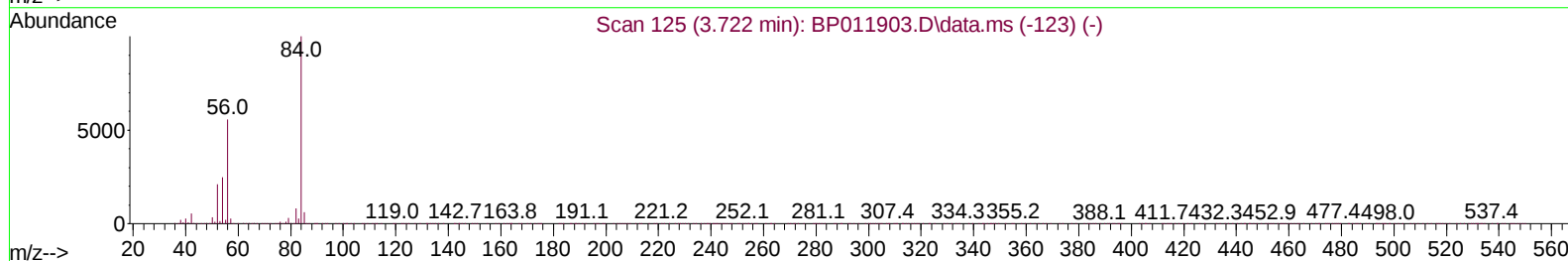
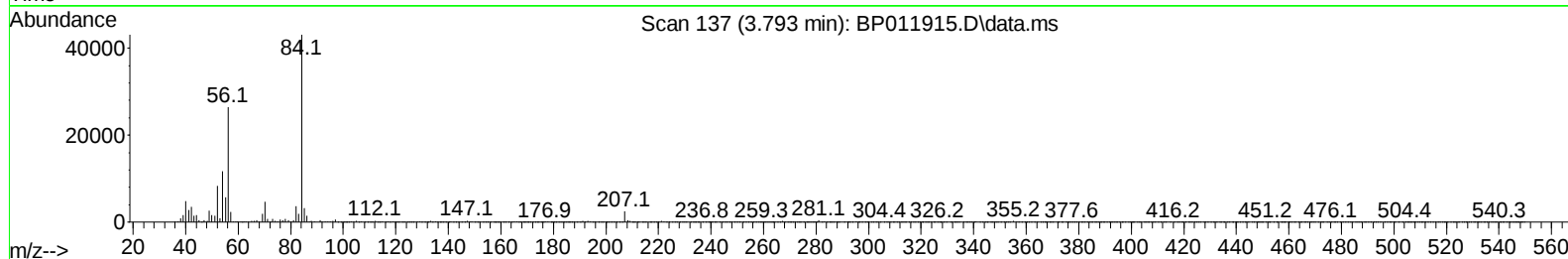
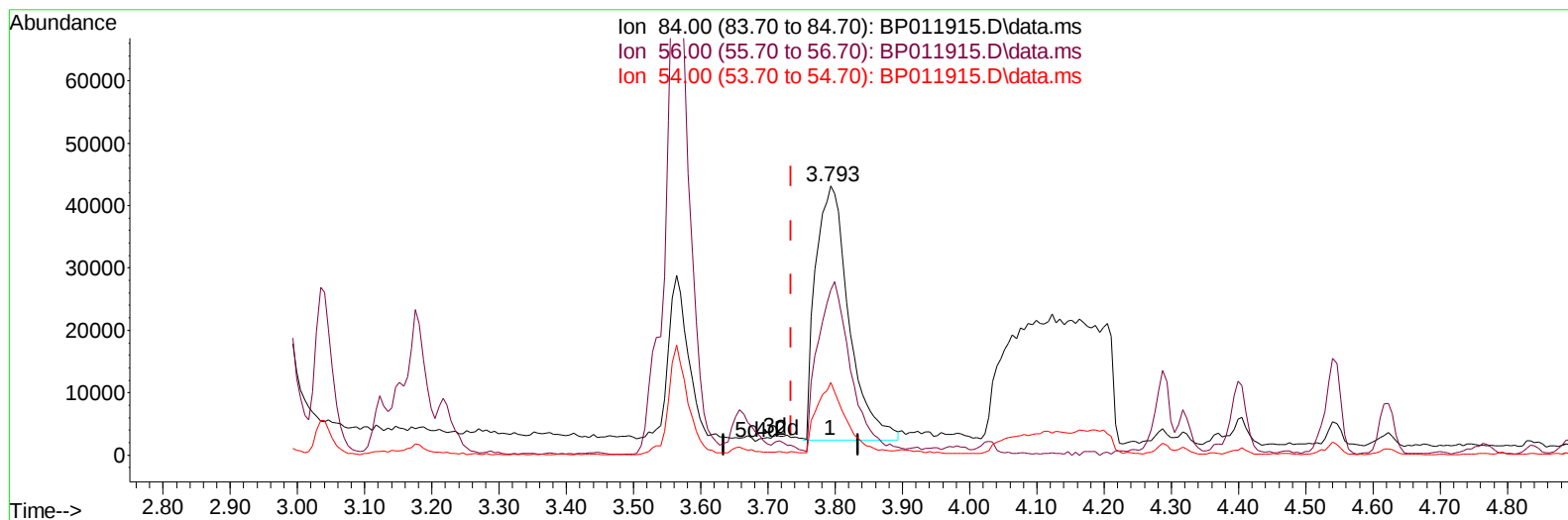
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TIC: BP011915.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.059) 9.39 ng/u1 m

response 140092

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	56.10	61.35
54.00	24.80	26.96
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	7.887	152	224598	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	902837	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.534	164	542359	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1104135	20.000	ng/ul	-0.01
79) Chrysene-d12	21.375	240	1083730	20.000	ng/ul	-0.01
88) Perylene-d12	23.810	264	1130268	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	29206	5.735	ng/uL	0.00
4) Pyridine-d5	3.793	84	140092m	9.392	ng/ul	0.06
7) Phenol-d5	7.058	99	122217	6.303	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	354340	32.969	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	393010	25.677	ng/ul	0.00
15) 4-Methylphenol-d8	8.599	113	231395	14.437	ng/ul	0.00
21) Nitrobenzene-d5	9.052	128	296978	43.185	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.775	143	266736	36.254	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	433815	31.878	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	438821	21.151	ng/ul	-0.01
46) Dimethylphthalate-d6	13.940	166	1525330	39.328	ng/ul	-0.01
49) Acenaphthylene-d8	14.222	160	1753967	39.670	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.740	143	45949	5.721	ng/ul	0.00
60) Fluorene-d10	15.528	176	1361572	40.424	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	194877	29.160	ng/ul	-0.01
73) Anthracene-d10	17.386	188	2131372	42.377	ng/ul	-0.01
81) Pyrene-d10	19.622	212	2510785	45.746	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2533698	44.781	ng/ul	-0.01

Target Compounds Qvalue

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

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