

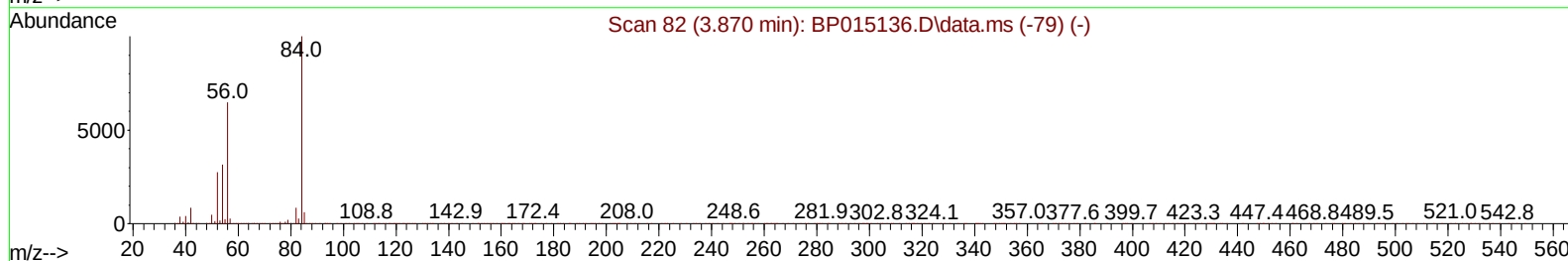
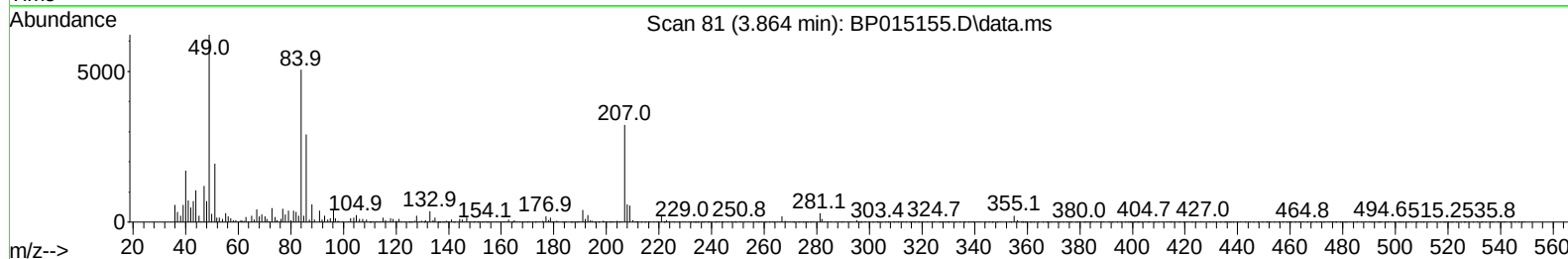
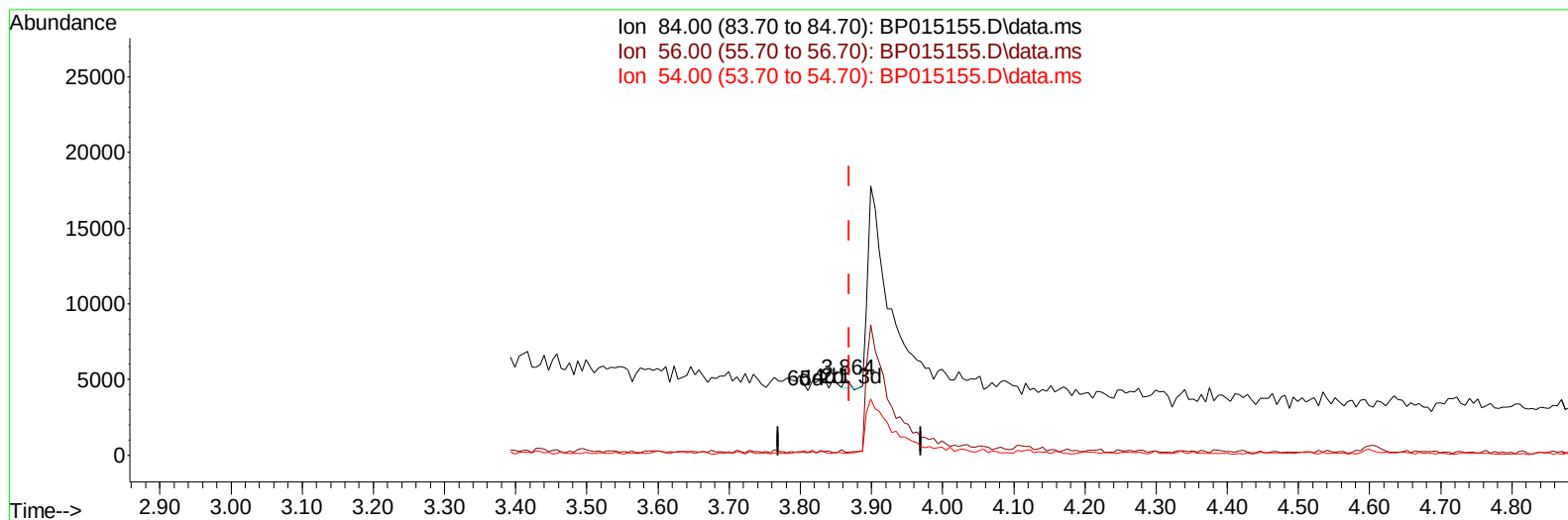
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051823\
Data File : BP015155.D
Acq On : 18 May 2023 20:35
Operator : CG/JU
Sample : 02612-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREG8

Manual IntegrationsAPPROVED

Quant Time: May 19 01:10:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 19 01:07:51 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/19/2023
Supervised By :Jagrut Upadhyay 05/19/2023



TIC: BP015155.D\data.ms

(4) Pyridine-d5 (S)

3.864min (-0.006) 0.01 ng/u/l

response 272

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	3.94#
54.00	29.10	2.10#
0.00	0.00	0.00

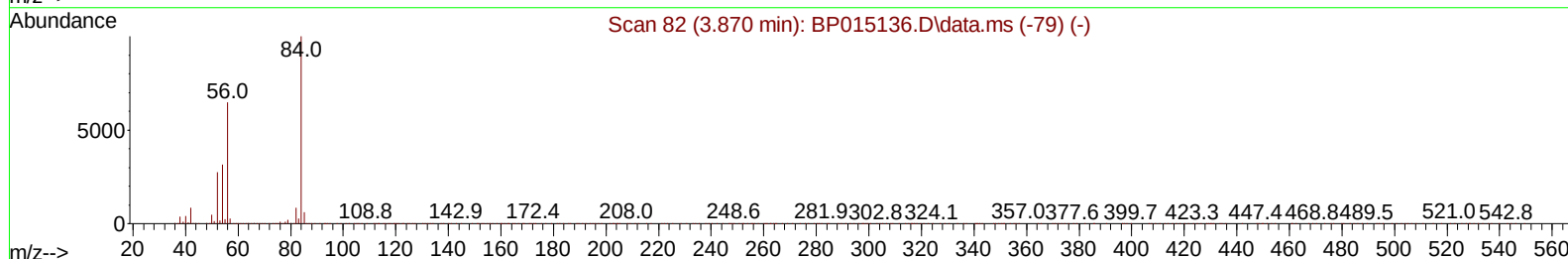
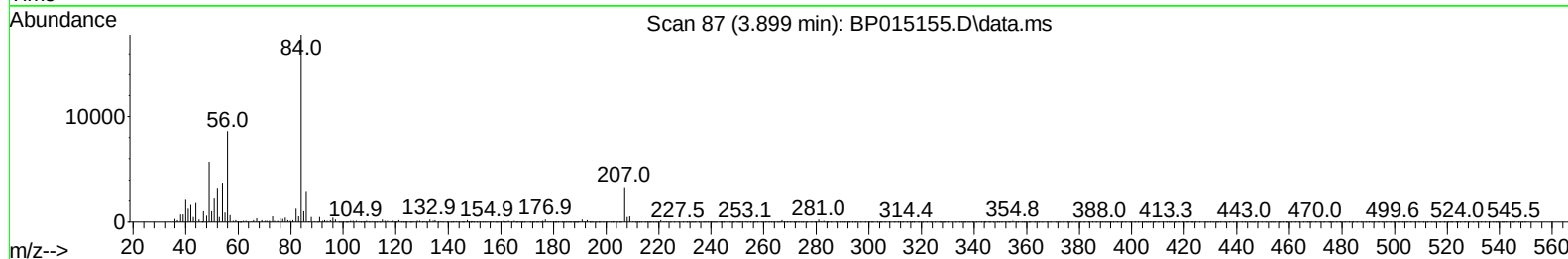
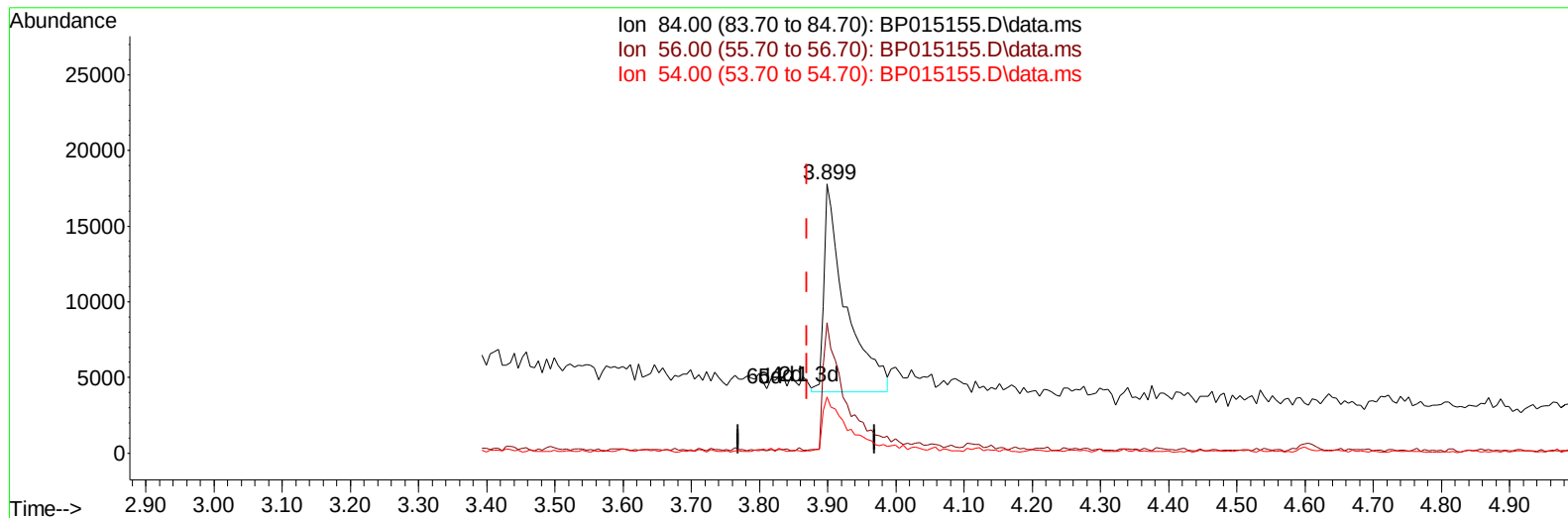
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TIC: BP015155.D\data.ms

(4) Pyridine-d5 (S)

3.899min (+ 0.029) 1.40 ng/ul m

response 30447

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	48.44#
54.00	29.10	20.93#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/19/2023
 Supervised By :Jagrut Upadhyay 05/19/2023

Quant Time: May 19 01:31:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 19 01:07:51 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.128	152	298202	20.000	ng/ul	0.00
20) Naphthalene-d8	10.957	136	1345366	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.763	164	847336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.516	188	1886505	20.000	ng/ul	0.00
79) Chrysene-d12	21.592	240	1652016	20.000	ng/ul	0.00
88) Perylene-d12	24.151	264	1717045	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	17812	2.321	ng/uL	0.00
4) Pyridine-d5	3.899	84	30447m	1.400	ng/ul	0.03
7) Phenol-d5	7.275	99	489796	17.690	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	340756	21.309	ng/ul	0.00
11) 2-Chlorophenol-d4	7.652	132	405879	19.938	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	420002	19.037	ng/ul	0.00
21) Nitrobenzene-d5	9.305	128	204466	19.341	ng/ul	0.00
24) 2-Nitrophenol-d4	10.028	143	221075	18.794	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	398071	18.754	ng/ul	0.00
31) 4-Chloroaniline-d4	11.093	131	563469	17.942	ng/ul	0.00
46) Dimethylphthalate-d6	14.169	166	1482901	23.160	ng/ul	0.00
49) Acenaphthylene-d8	14.463	160	1633411	22.305	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	177262	14.257	ng/ul	0.00
60) Fluorene-d10	15.757	176	1262238	23.362	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	148792	12.515	ng/ul	0.00
73) Anthracene-d10	17.616	188	2017379	23.354	ng/ul	0.00
81) Pyrene-d10	19.833	212	2404539	25.096	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.986	264	2150764	25.233	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.963	105	79720	2.315	ng/ul	96

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

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