

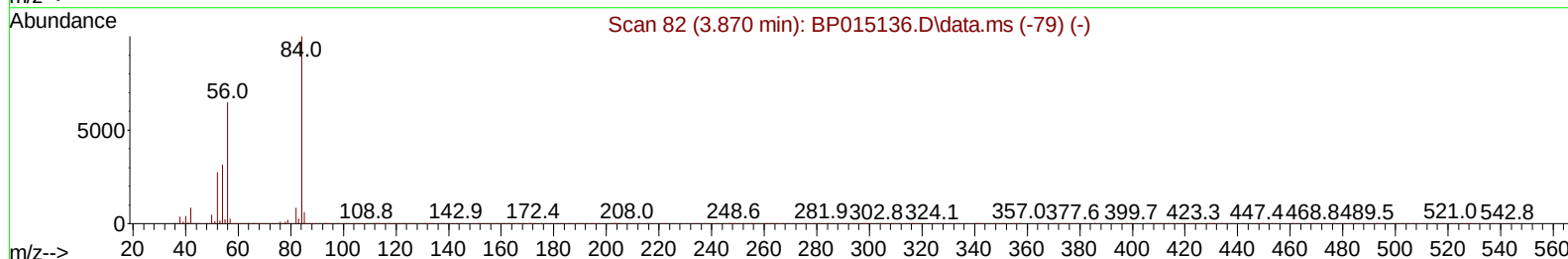
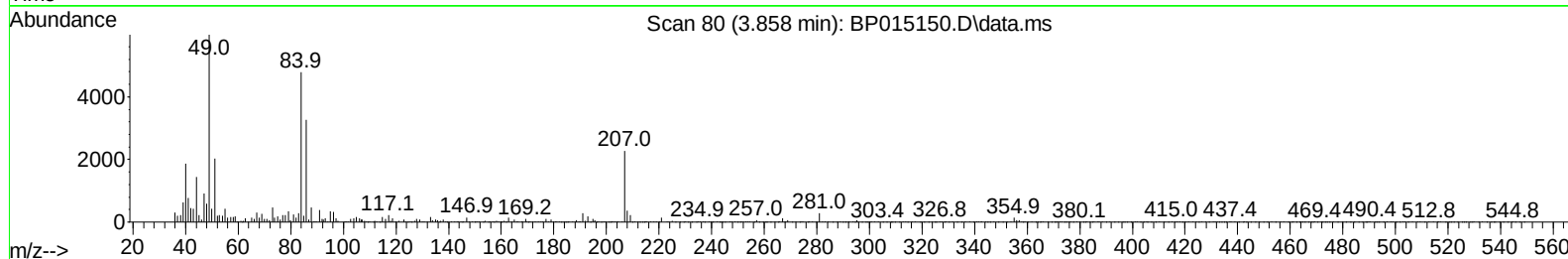
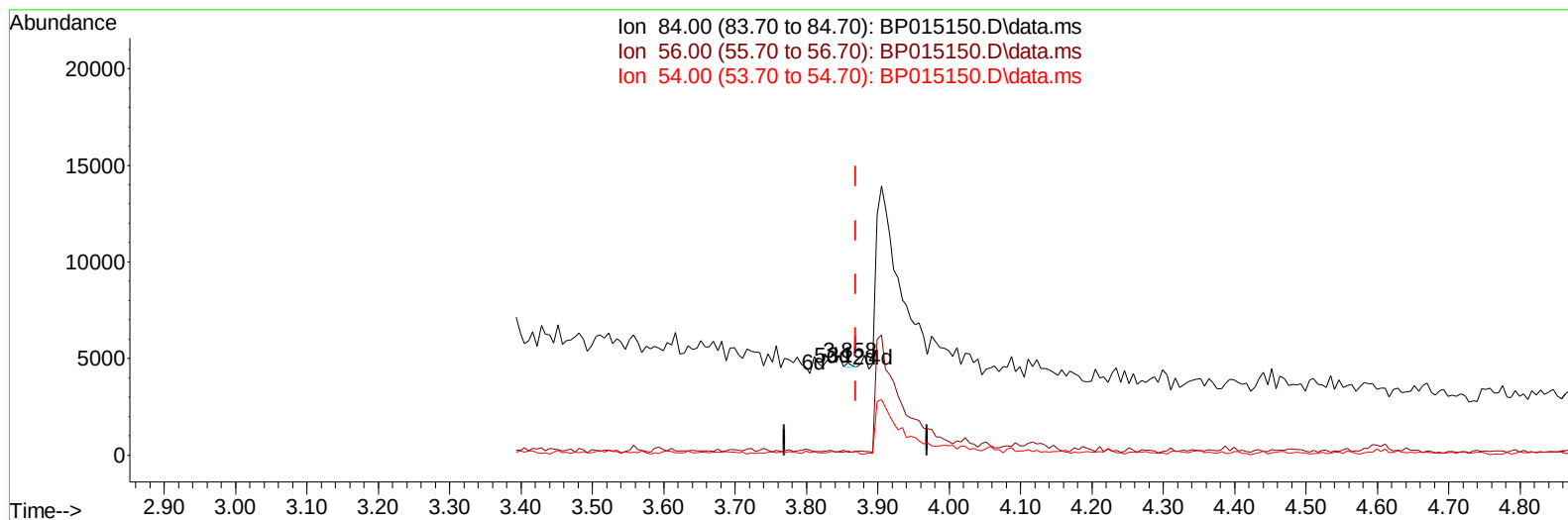
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051823\
 Data File : BP015150.D
 Acq On : 18 May 2023 17:48
 Operator : CG/JU
 Sample : 02612-21
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREB0

Manual IntegrationsAPPROVED

Quant Time: May 19 01:09:43 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: BP015150.D\data.ms

(4) Pyridine-d5 (S)

3.858min (-0.012) 0.01 ng/uI

response 115

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	3.16#
54.00	29.10	4.29#
0.00	0.00	0.00

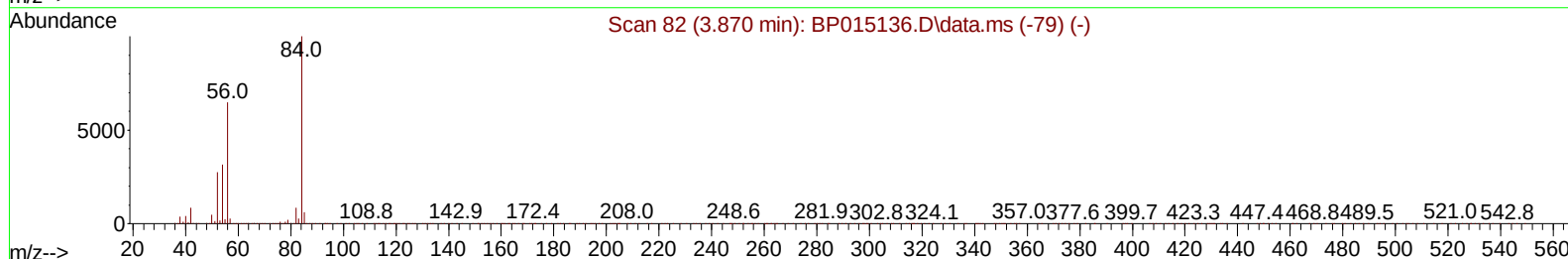
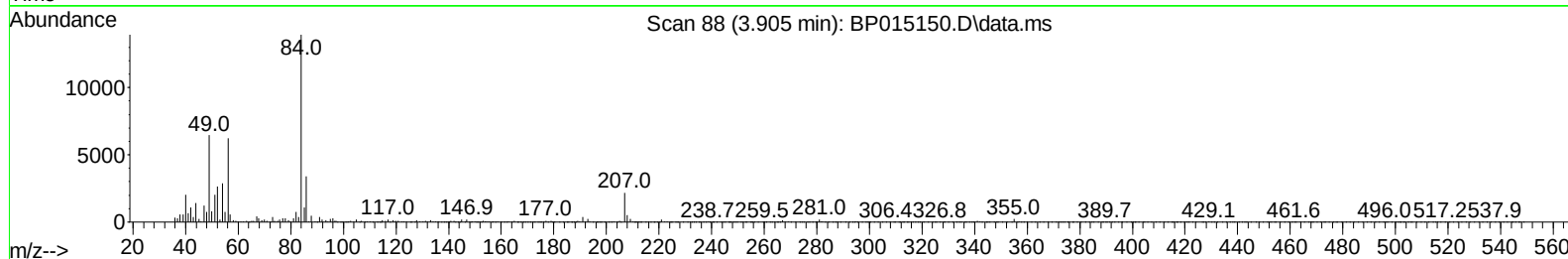
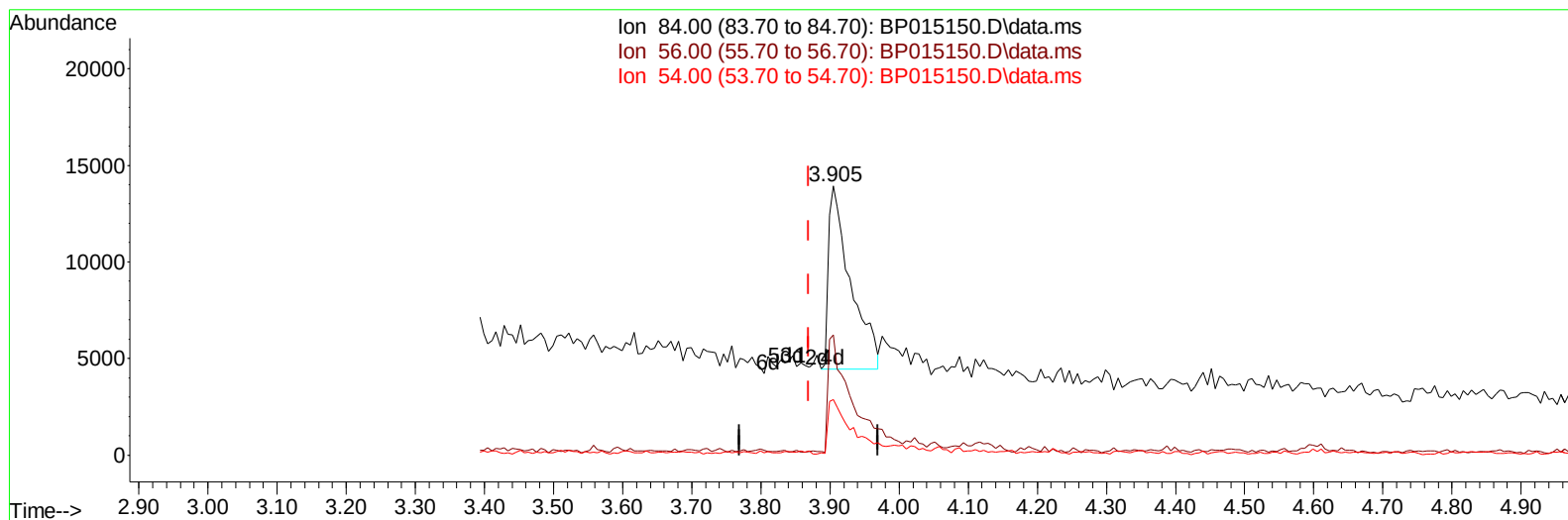
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TIC: BP015150.D\data.ms

(4) Pyridine-d5 (S)

3.905min (+ 0.035) 1.03 ng/ul m

response 21098

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.60	44.76#
54.00	29.10	20.54#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051823\
 Data File : BP051510.D
 Acq On : 18 May 2023 17:48
 Operator : CG/JU
 Sample : 02612-21
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREB0

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/19/2023
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Quant Time: May 19 01:24:12 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	280408	20.000	ng/ul	0.00
20) Naphthalene-d8	10.957	136	1279812	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.769	164	820157	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.516	188	1810330	20.000	ng/ul	0.00
79) Chrysene-d12	21.592	240	1595980	20.000	ng/ul	0.00
88) Perylene-d12	24.151	264	1687729	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	16963	2.350	ng/uL	0.00
4) Pyridine-d5	3.905	84	21098m	1.031	ng/ul	0.04
7) Phenol-d5	7.275	99	528311	20.292	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.446	67	349348	23.233	ng/ul	0.00
11) 2-Chlorophenol-d4	7.652	132	410055	21.421	ng/ul	0.00
15) 4-Methylphenol-d8	8.834	113	423527	20.415	ng/ul	0.00
21) Nitrobenzene-d5	9.305	128	215552	21.434	ng/ul	0.00
24) 2-Nitrophenol-d4	10.034	143	237289	21.206	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	422019	20.900	ng/ul	0.00
31) 4-Chloroaniline-d4	11.093	131	606341	20.296	ng/ul	0.00
46) Dimethylphthalate-d6	14.169	166	1540285	24.853	ng/ul	0.00
49) Acenaphthylene-d8	14.463	160	1707508	24.089	ng/ul	0.00
54) 4-Nitrophenol-d4	14.951	143	211211	17.551	ng/ul	0.00
60) Fluorene-d10	15.757	176	1324912	25.335	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	171721	15.052	ng/ul	0.00
73) Anthracene-d10	17.616	188	2109204	25.444	ng/ul	0.00
81) Pyrene-d10	19.833	212	2757365	29.789	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.986	264	2413888	28.812	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.963	105	130285	4.024	ng/ul	96

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

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