

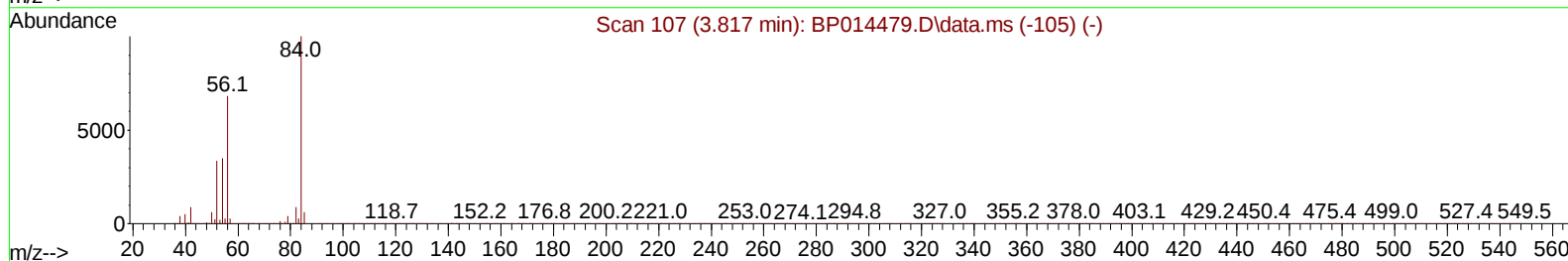
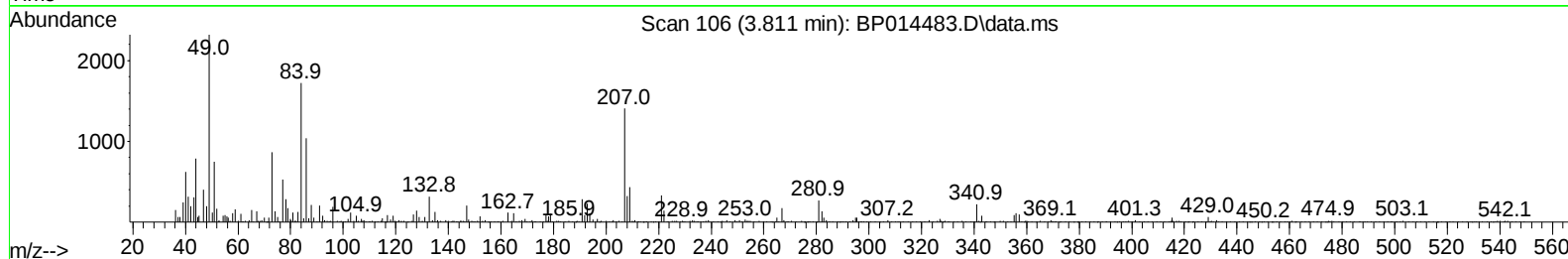
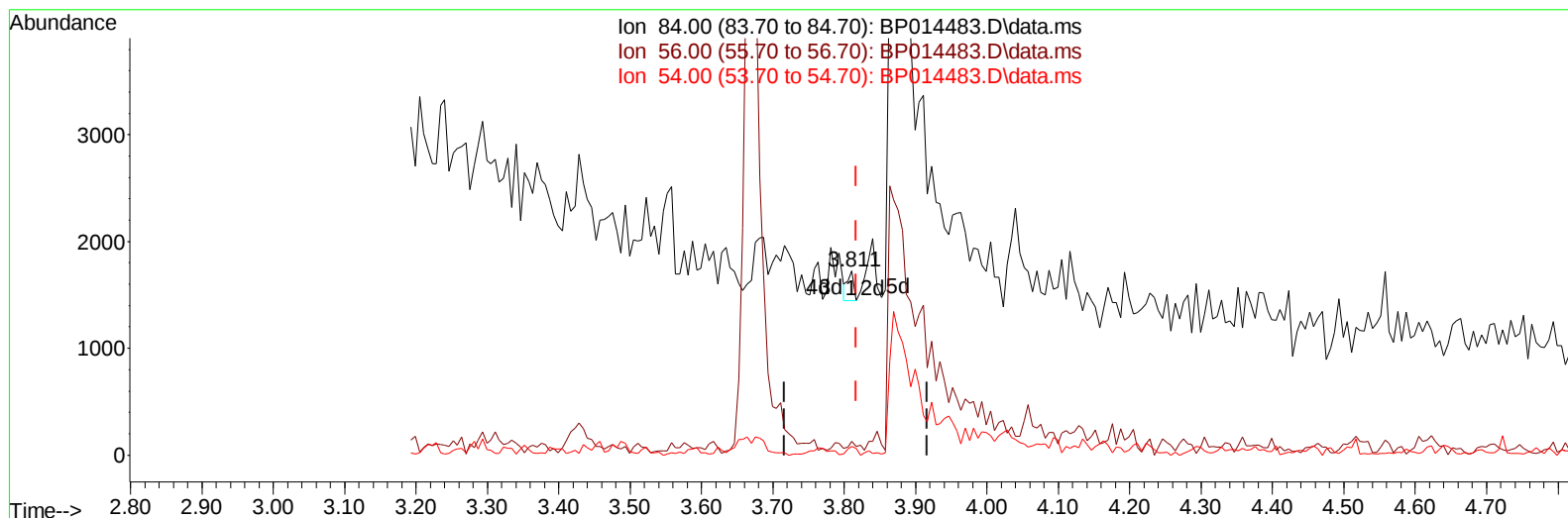
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014483.D
 Acq On : 03 Apr 2023 12:59
 Operator : CG/JU
 Sample : 02093-05DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB973DL

Manual IntegrationsAPPROVED

Quant Time: Apr 04 00:52:10 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 04/04/2023
 Supervised By :Jagrut Upadhyay 04/04/2023



TIC: BP014483.D\data.ms

(4) Pyridine-d5 (S)

3.811min (-0.006) 0.01 ng/ul

response 163

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	7.48#
54.00	32.60	4.58#
0.00	0.00	0.00

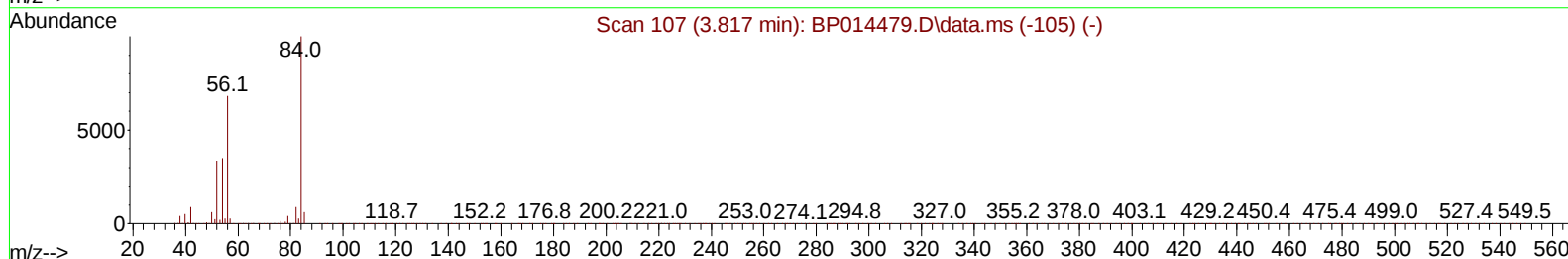
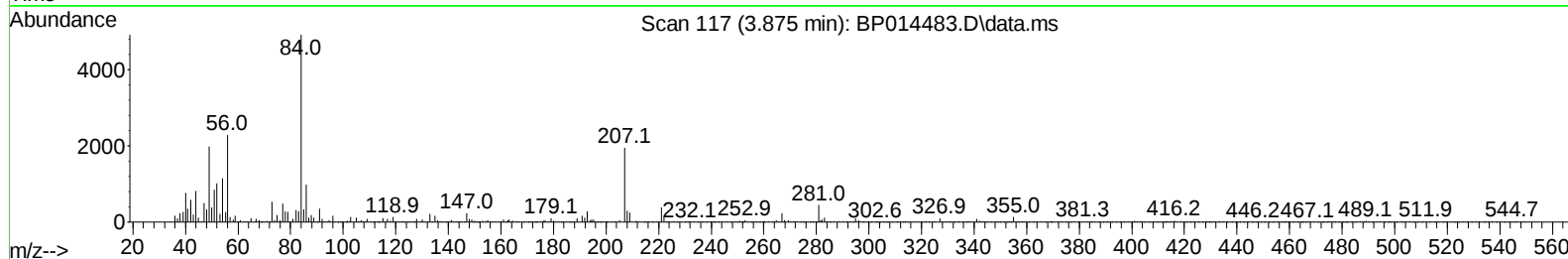
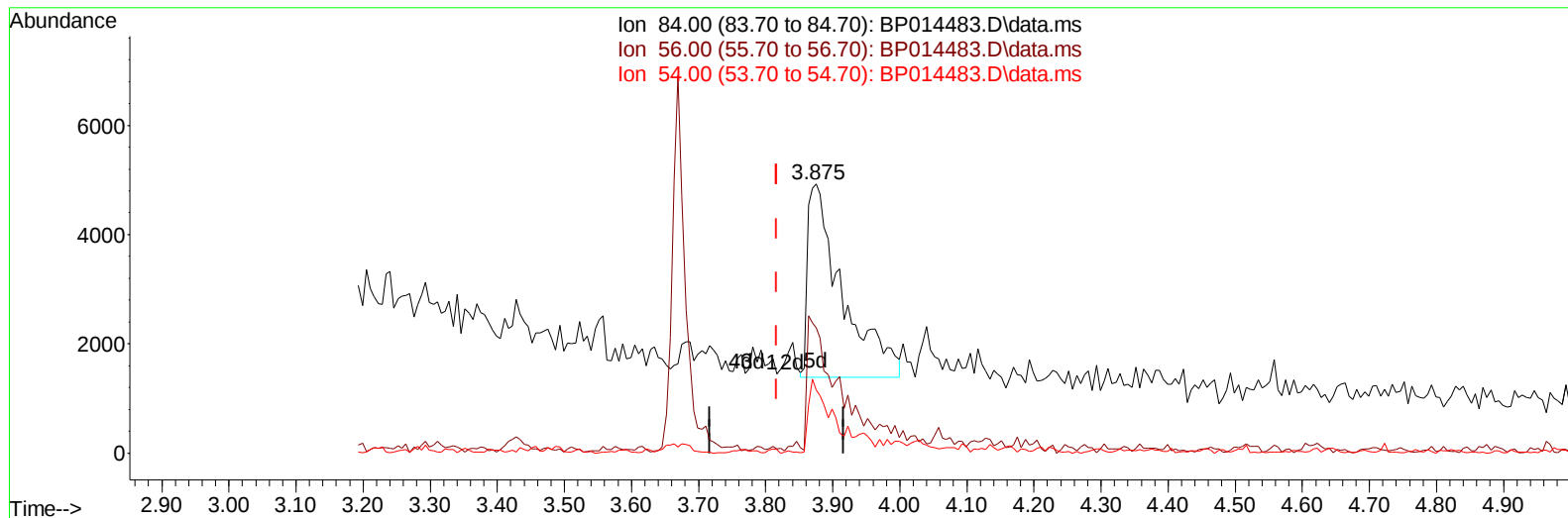
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TIC: BP014483.D\data.ms

(4) Pyridine-d5 (S)

3.875min (+ 0.059) 1.11 ng/ul m

response 12666

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	46.62#
54.00	32.60	23.48#
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	173407	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	779464	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.663	164	534571	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	1165005	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	879817	20.000	ng/ul	0.00
88) Perylene-d12	24.033	264	849592	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	1895	0.424	ng/uL	0.00
4) Pyridine-d5	3.875	84	12666m	1.105	ng/ul	0.06
7) Phenol-d5	7.199	99	12769	0.798	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.334	67	33402	3.255	ng/ul	0.00
11) 2-Chlorophenol-d4	7.546	132	31708	2.743	ng/ul	0.00
15) 4-Methylphenol-d8	8.746	113	23827	1.847	ng/ul	0.02
21) Nitrobenzene-d5	9.193	128	18009	2.862	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	20597	2.859	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	34768	2.650	ng/ul	0.02
31) 4-Chloroaniline-d4	10.993	131	51317	2.969	ng/ul	0.01
46) Dimethylphthalate-d6	14.069	166	161509	3.726	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	164857	3.428	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.657	176	136951	3.789	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.528	188	208384	3.696	ng/ul	0.00
81) Pyrene-d10	19.757	212	238070	4.028	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.863	264	162441	3.599	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.422	88	79892	16.258	ng/uL	96

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

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