

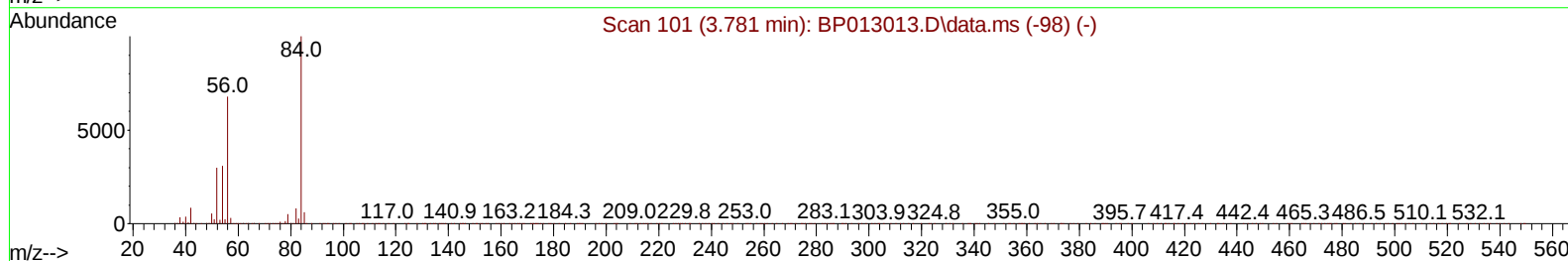
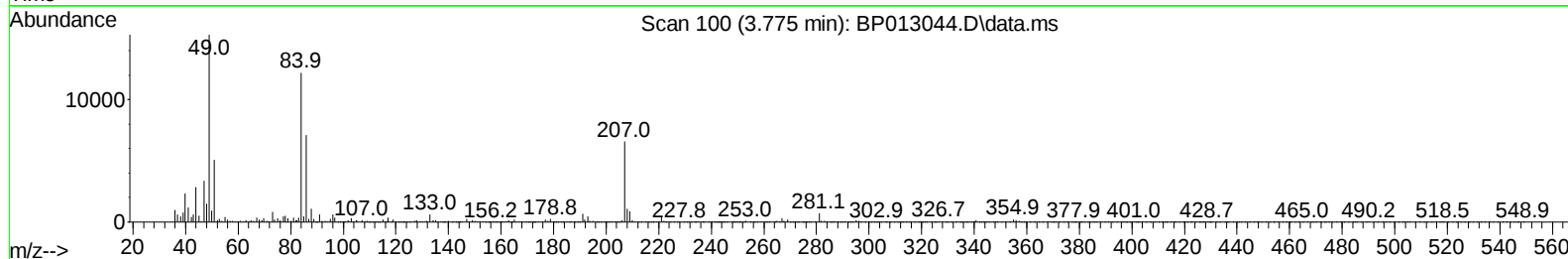
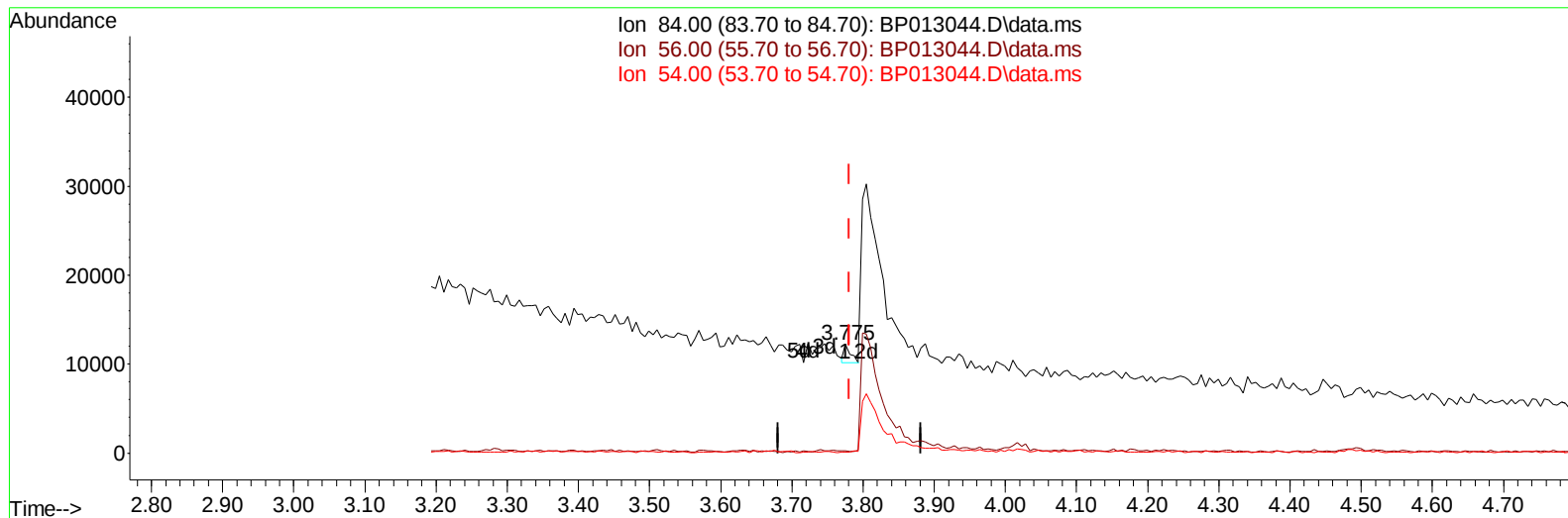
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013044.D
 Acq On : 10 Dec 2022 17:35
 Operator : CG/JU
 Sample : N5950-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYF49

Manual IntegrationsAPPROVED

Quant Time: Dec 12 03:25:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/12/2022
 Supervised By :mohammad ahmed 12/12/2022



TIC: BP013044.D\data.ms

(4) Pyridine-d5 (S)

3.775min (-0.006) 0.03 ng/u1

response 1369

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	1.91#
54.00	30.50	0.92#
0.00	0.00	0.00

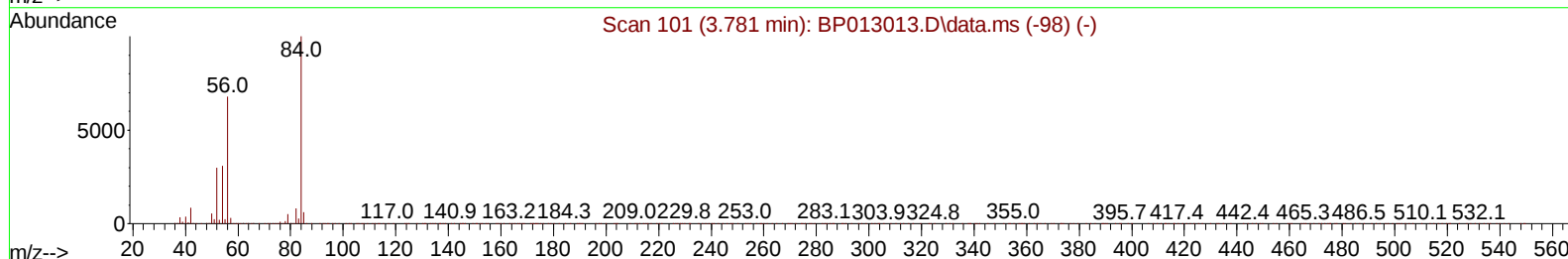
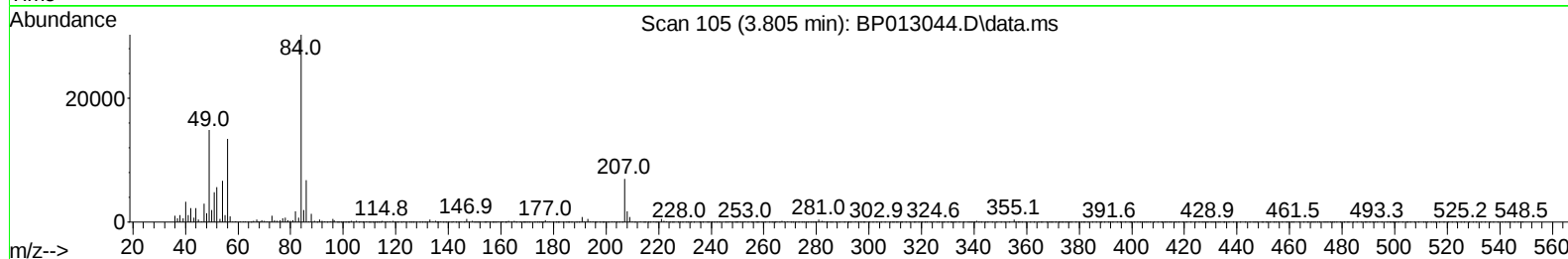
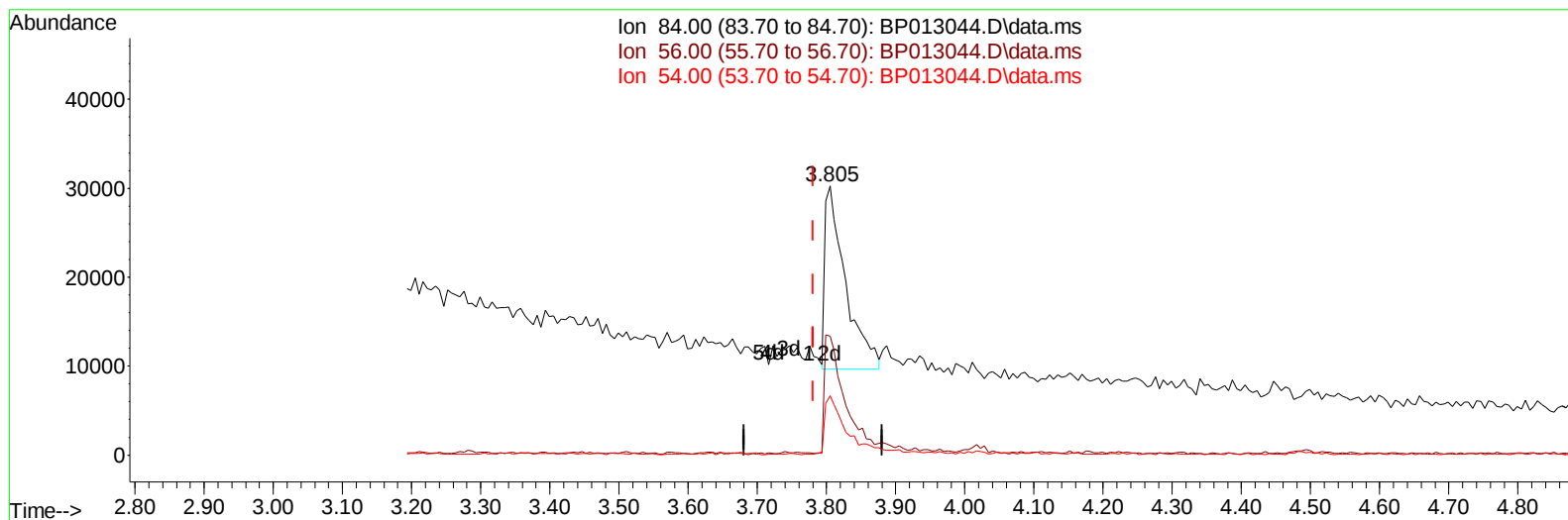
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TIC: BP013044.D\data.ms

(4) Pyridine-d5 (S)

3.805min (+ 0.024) 0.93 ng/u1 m

response 42617

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	44.25#
54.00	30.50	21.99#
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	7.969	152	691432	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	3060475	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	2076653	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	4539160	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3413655	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	3101832	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	73170	4.525	ng/uL	0.00
4) Pyridine-d5	3.805	84	42617m	0.928	ng/ul	0.02
7) Phenol-d5	7.128	99	375584	6.669	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	1076095	31.674	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	1010825	22.800	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	676603	14.662	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	750285	33.367	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	795016	31.405	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1318606	26.817	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	374225	5.391	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	5447280	34.131	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	5922466	33.771	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	141443	4.927	ng/ul	0.00
60) Fluorene-d10	15.604	176	4726876	35.008	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	877460	30.039	ng/ul	0.00
73) Anthracene-d10	17.463	188	7821259	38.193	ng/ul	0.00
81) Pyrene-d10	19.698	212	8705634	44.429	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.774	264	6197793	40.087	ng/ul	0.00

Target Compounds Qvalue

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

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