

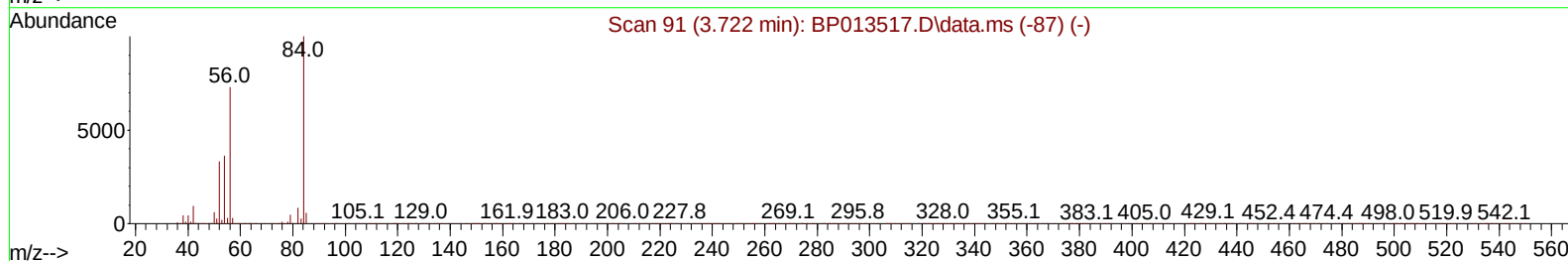
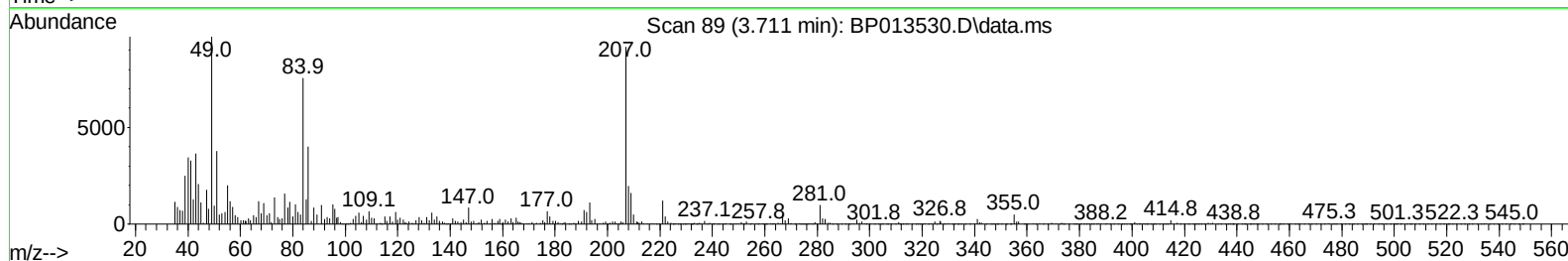
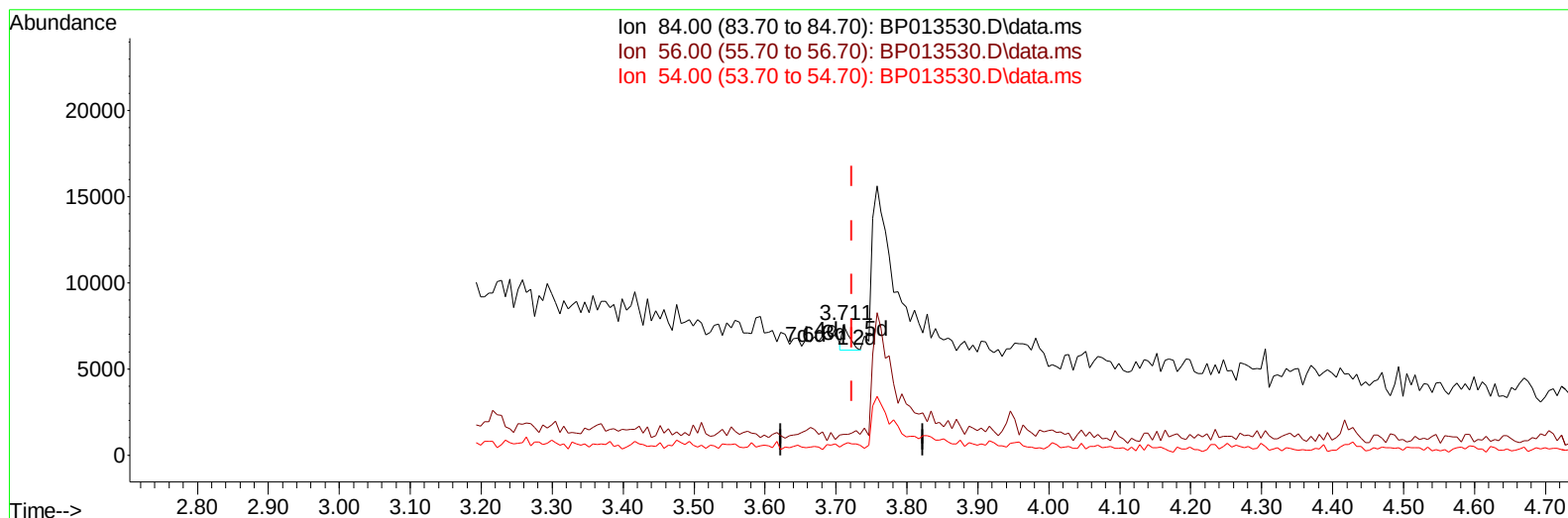
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011823\
Data File : BP013530.D
Acq On : 18 Jan 2023 19:55
Operator : CG/JU
Sample : 01146-04
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Z9

Manual IntegrationsAPPROVED

Quant Time: Jan 19 01:48:49 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010523.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 18 01:06:35 2023
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/19/2023
Supervised By :mohammad ahmed 01/20/2023



TIC: BP013530.D\data.ms

(4) Pyridine-d5 (S)

3.711min (-0.012) 0.04 ng/uI

response 1061

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.30	15.70#
54.00	34.30	8.30#
0.00	0.00	0.00

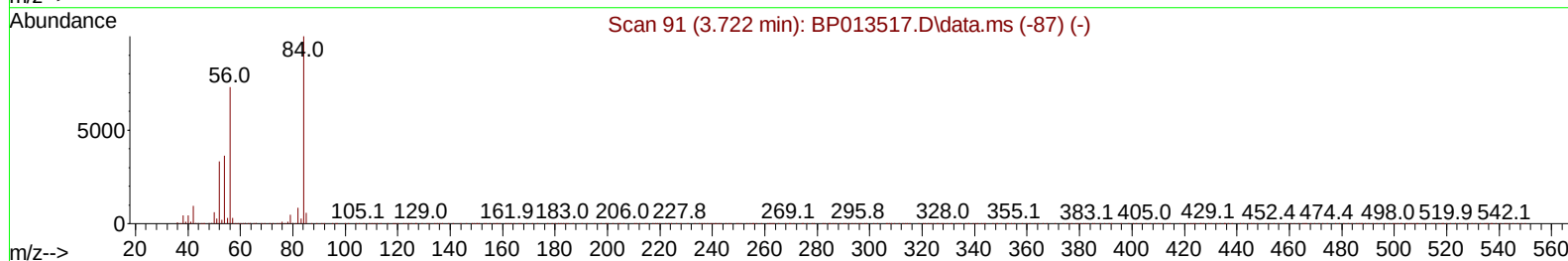
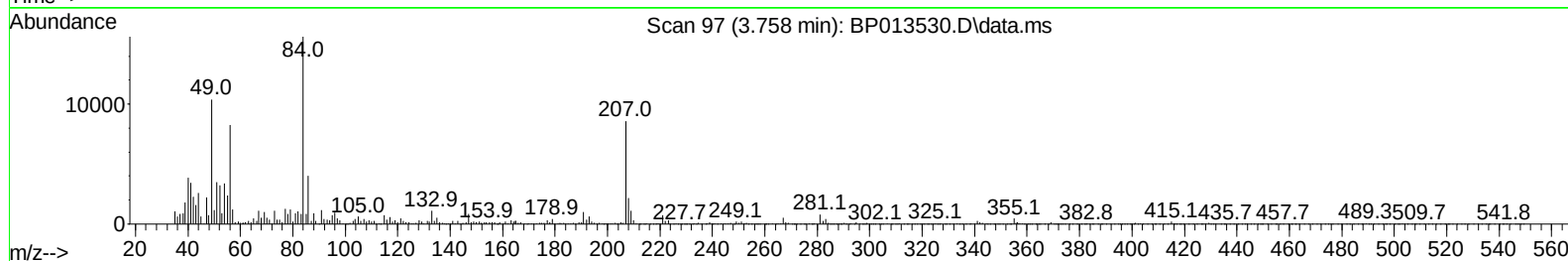
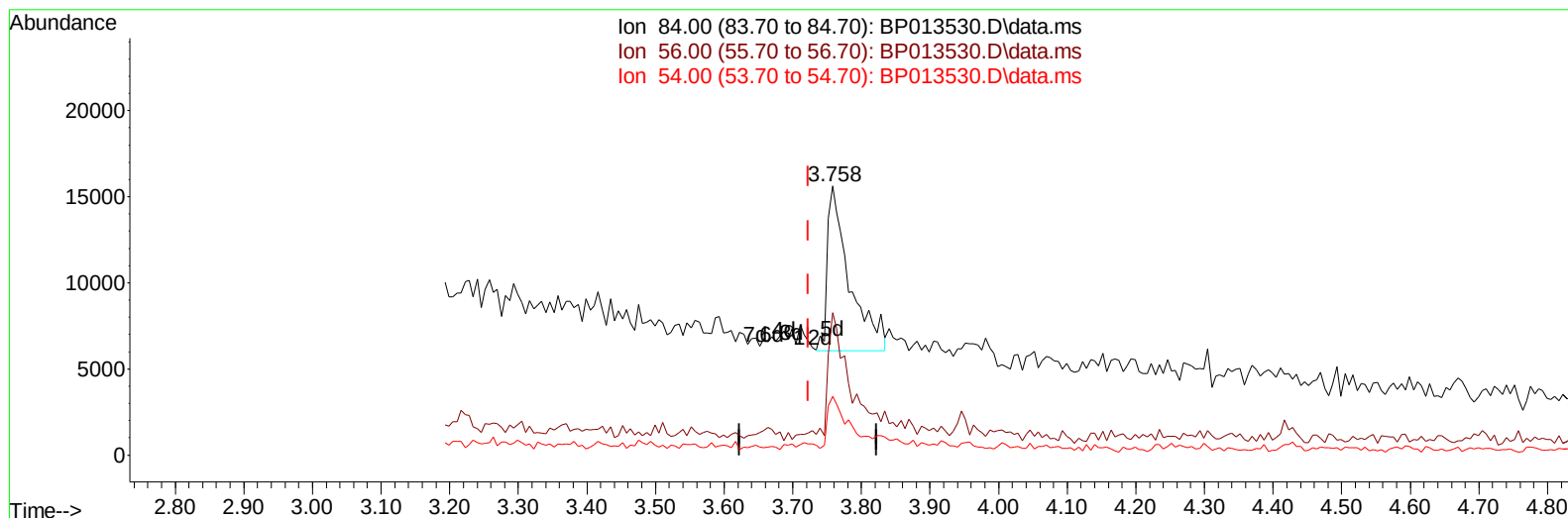
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TIC: BP013530.D\data.ms

(4) Pyridine-d5 (S)

3.758min (+ 0.035) 0.73 ng/ul m

response 21565

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	72.30	52.85#
54.00	34.30	21.76#
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.899	152		411828	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136		1608902	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164		935878	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.304	188		1871186	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240		1772400	20.000	ng/ul	0.00
88) Perylene-d12	23.839	264		2295141	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.293	96		40506	3.823	ng/uL	0.00
4) Pyridine-d5	3.758	84		21565m	0.730	ng/ul	0.04
7) Phenol-d5	7.058	99		751991	22.005	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67		495021	22.961	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132		623749	24.476	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113		593770	22.806	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128		297186	25.942	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143		334297	26.658	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165		596371	23.863	ng/ul	0.00
31) 4-Chloroaniline-d4	10.851	131		310911	8.831	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166		1710159	25.081	ng/ul	0.00
49) Acenaphthylene-d8	14.240	160		1965908	26.047	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143		164874	13.210	ng/ul	0.00
60) Fluorene-d10	15.539	176		1494933	24.822	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.663	200		199675	16.829	ng/ul	0.00
73) Anthracene-d10	17.404	188		2149048	25.984	ng/ul	0.00
81) Pyrene-d10	19.639	212		2624784	27.219	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264		3033831	26.910	ng/ul	-0.02
Target Compounds							
86) Bis(2-ethylhexyl)phtha...	21.286	149		2323609	47.007	ng/ul	99

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

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