

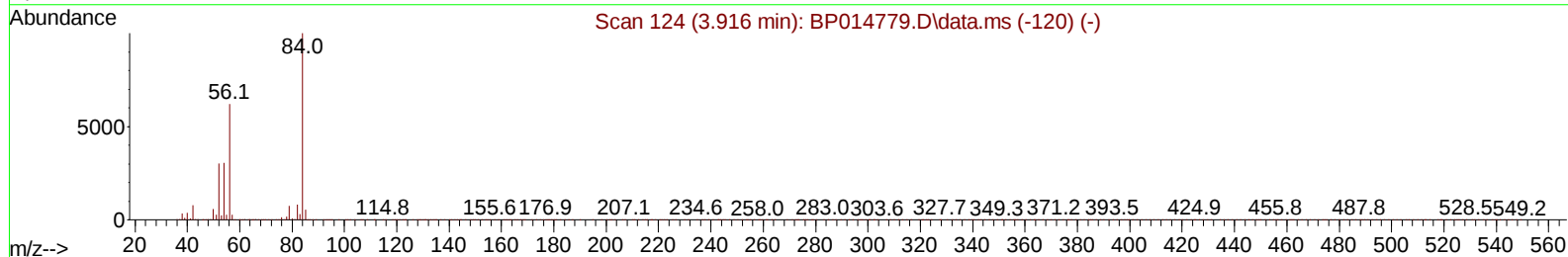
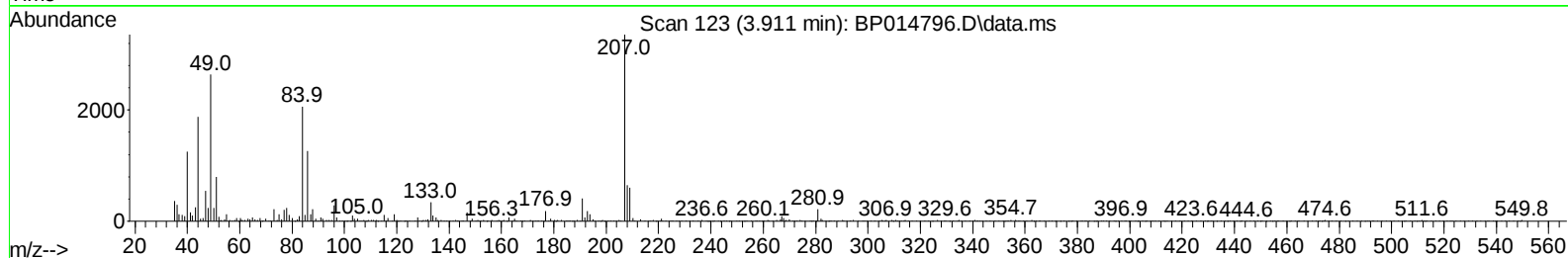
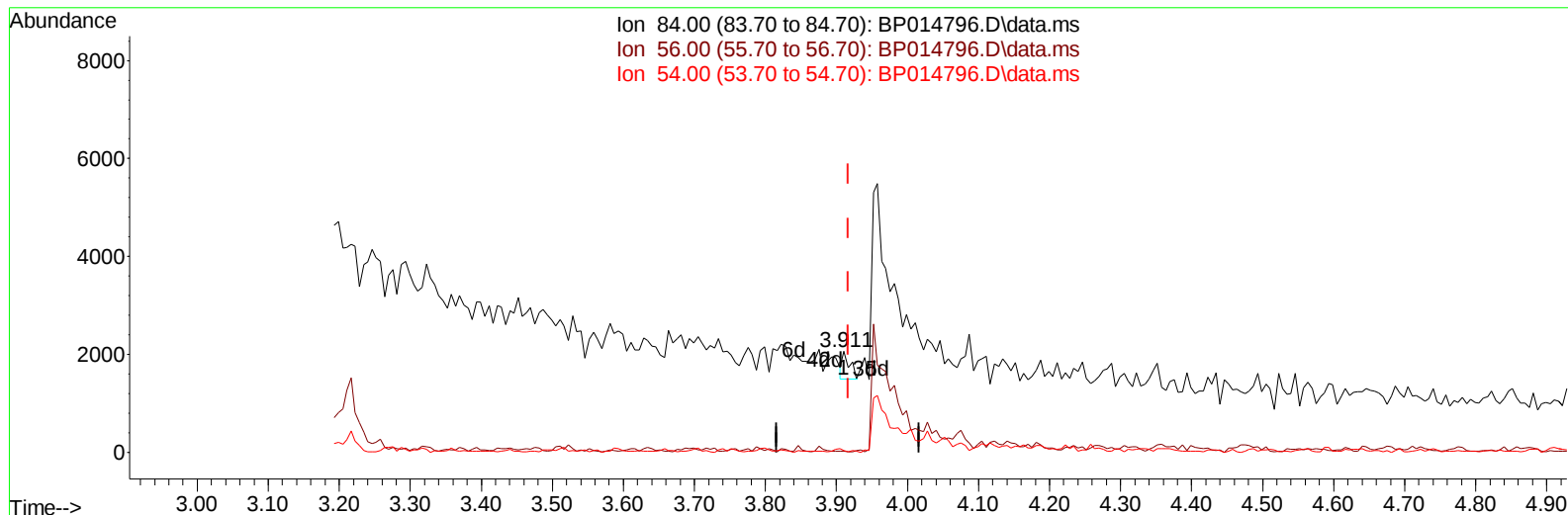
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
Data File : BP014796.D
Acq On : 21 Apr 2023 18:36
Operator : CG/JU
Sample : 02378-13DL 20X
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMK6DL

Manual IntegrationsAPPROVED

Quant Time: Apr 21 23:23:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 20 02:32:50 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023



TIC: BP014796.D\data.ms

(4) Pyridine-d5 (S)

3.911min (-0.006) 0.02 ng/u1

response 415

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.70	1.12#
54.00	30.50	1.89#
0.00	0.00	0.00

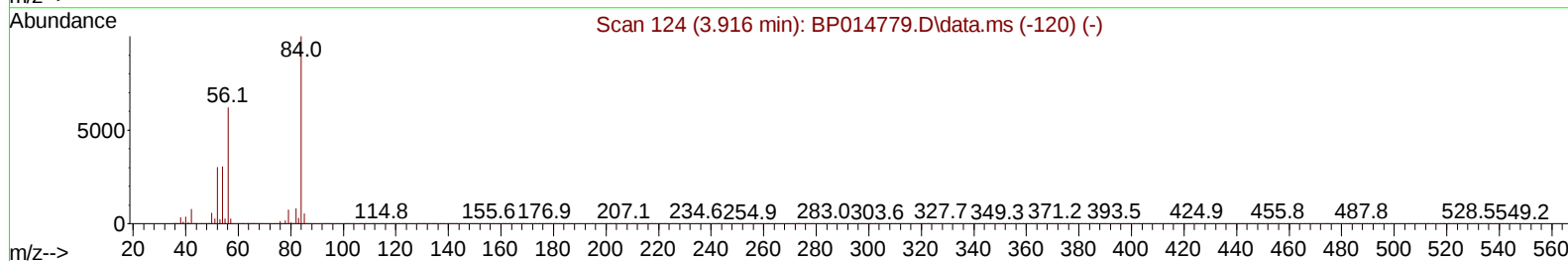
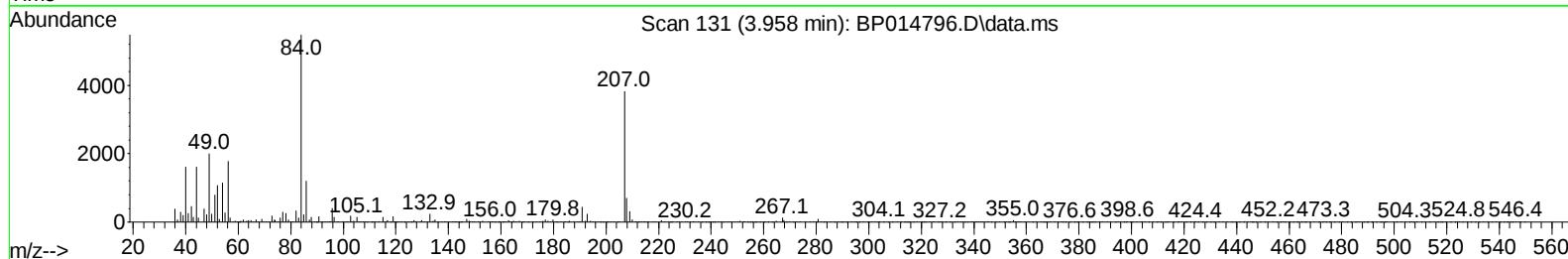
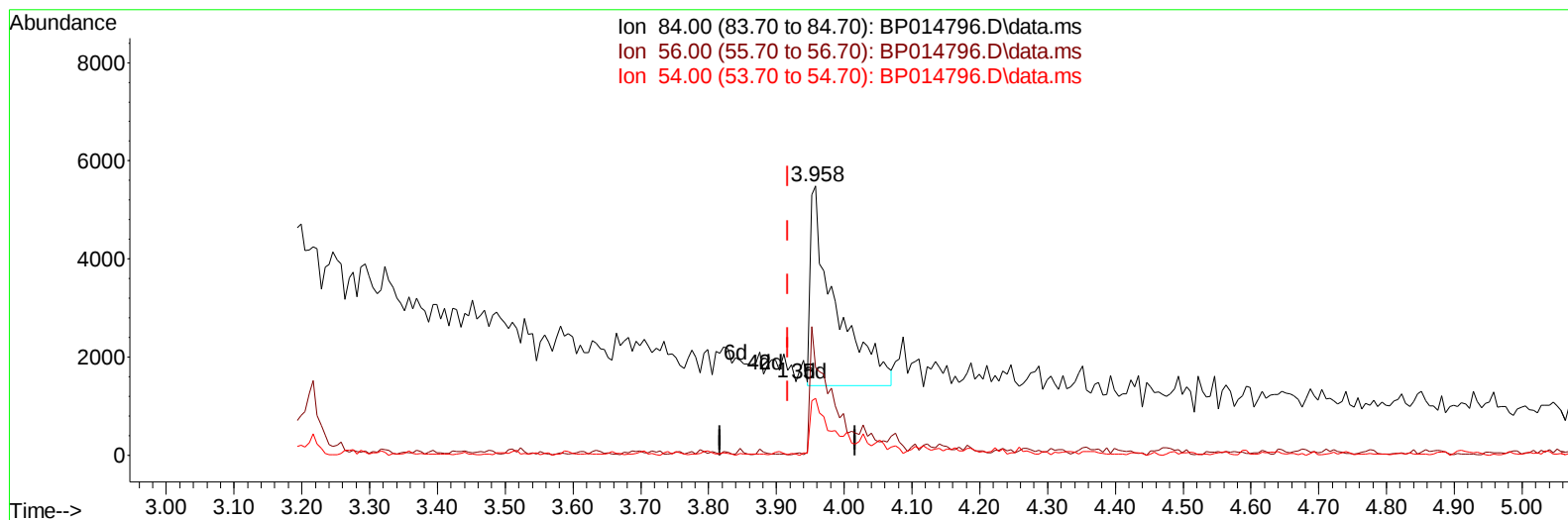
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TIC: BP014796.D\data.ms

(4) Pyridine-d5 (S)

3.958min (+ 0.041) 0.48 ng/ul m

response 10439

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.70	32.45#
54.00	30.50	21.19#
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	8.152	152	315471	20.000	ng/ul	0.00
20) Naphthalene-d8	10.981	136	1305334	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.781	164	841558	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.528	188	1808374	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	1211873	20.000	ng/ul	0.00
88) Perylene-d12	24.168	264	1151816	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.487	96	1935	0.244	ng/uL	0.00
4) Pyridine-d5	3.958	84	10439m	0.483	ng/ul	0.04
7) Phenol-d5	7.281	99	7992	0.293	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.463	67	27083	1.609	ng/ul	0.00
11) 2-Chlorophenol-d4	7.675	132	20320	1.003	ng/ul	0.00
15) 4-Methylphenol-d8	8.846	113	13761	0.647	ng/ul	0.00
21) Nitrobenzene-d5	9.322	128	11483	1.404	ng/ul	0.00
24) 2-Nitrophenol-d4	10.046	143	9755	1.320	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.587	165	20395	1.017	ng/ul	0.00
31) 4-Chloroaniline-d4	11.104	131	31300	1.022	ng/ul	0.00
46) Dimethylphthalate-d6	14.175	166	102495	1.538	ng/ul	0.00
49) Acenaphthylene-d8	14.475	160	96541	1.250	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.763	176	86876	1.519	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.628	188	128810	1.474	ng/ul	0.00
81) Pyrene-d10	19.839	212	133160	1.918	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.998	264	80621	1.350	ng/ul	-0.02
Target Compounds						
					Qvalue	
22) Nitrobenzene	9.363	77	979922	44.090	ng/ul	98

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

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