

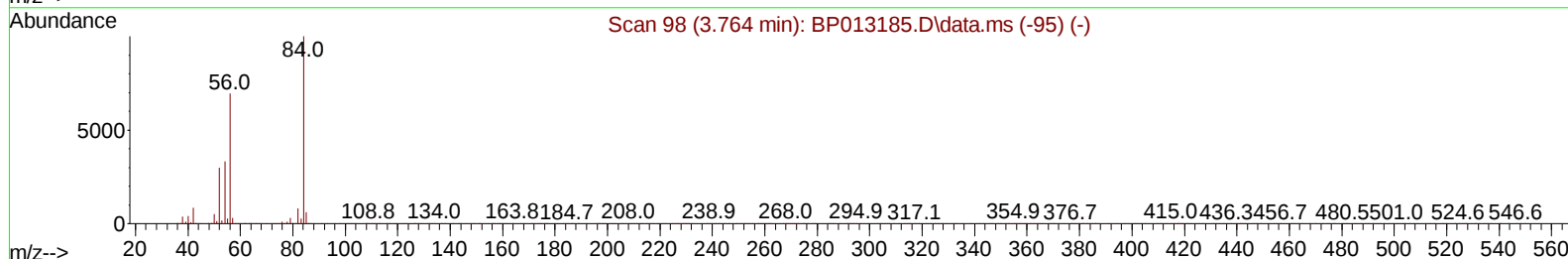
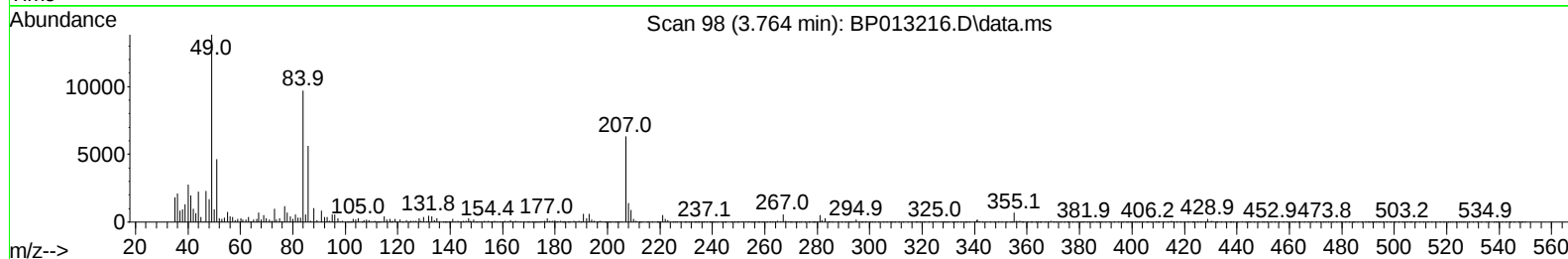
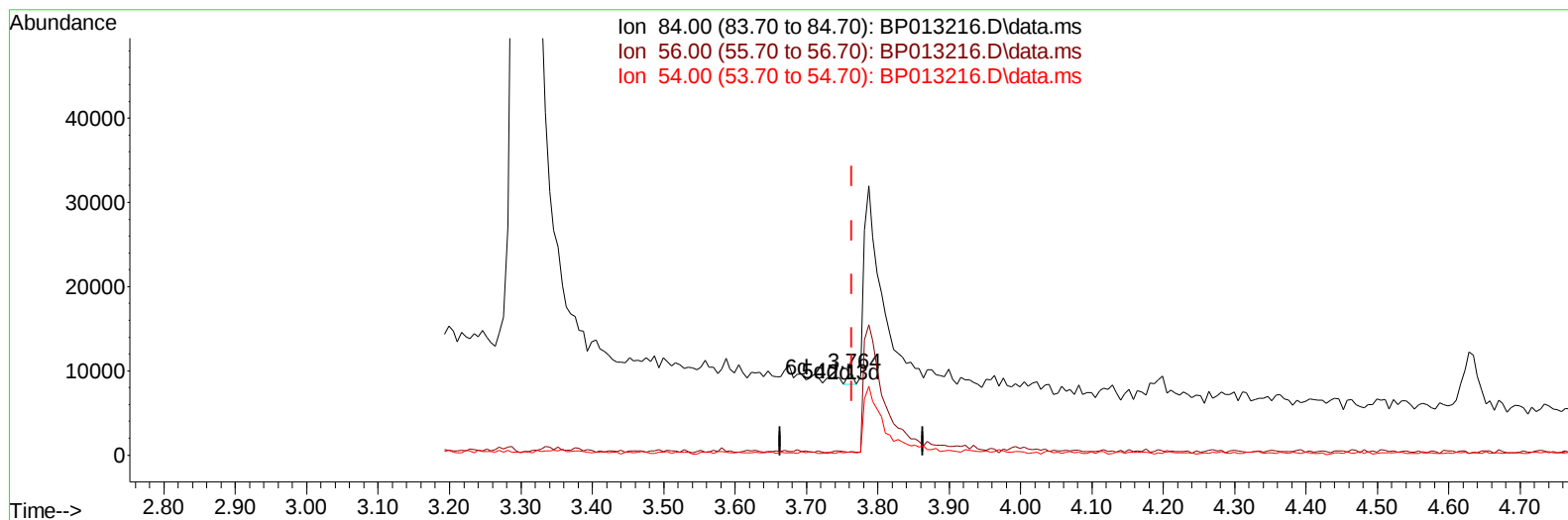
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
 Data File : BP013216.D
 Acq On : 21 Dec 2022 19:46
 Operator : CG/JU
 Sample : N6032-03
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COL86

Manual IntegrationsAPPROVED

Quant Time: Dec 21 21:55:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 19 21:43:11 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/22/2022
 Supervised By :mohammad ahmed 12/22/2022



TIC: BP013216.D\data.ms

(4) Pyridine-d5 (S)

3.764min (+ 0.000) 0.02 ng/u1

response 823

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	4.31#
54.00	30.50	3.24#
0.00	0.00	0.00

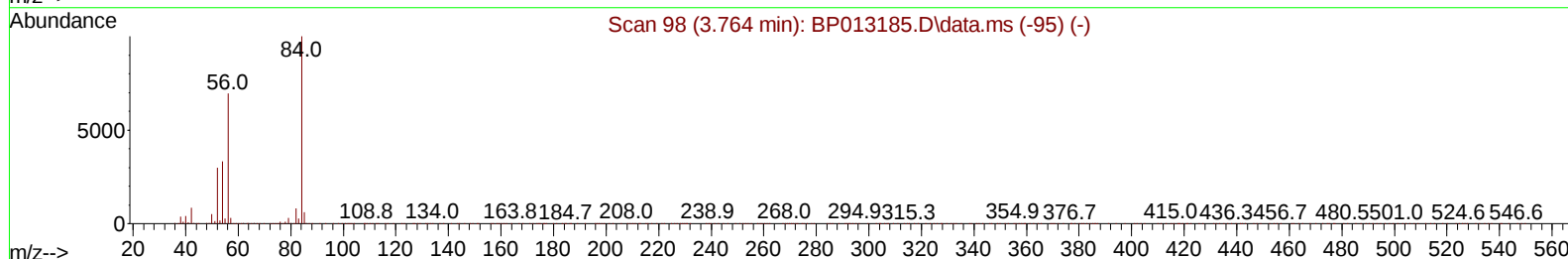
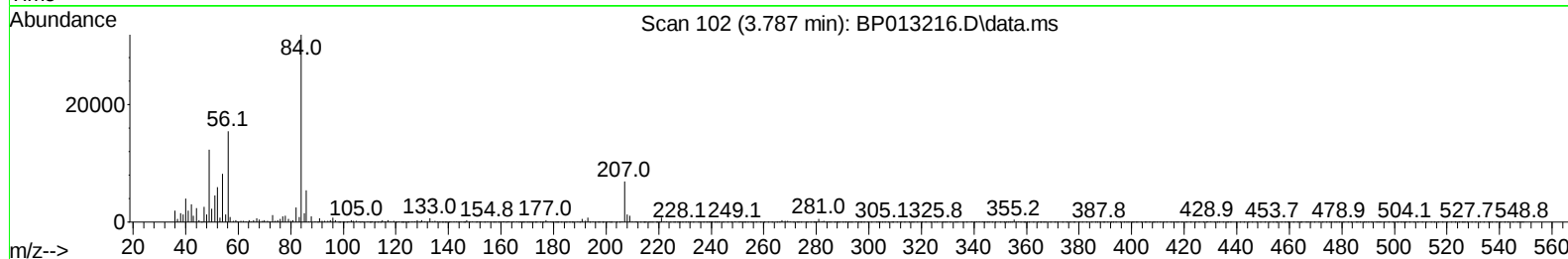
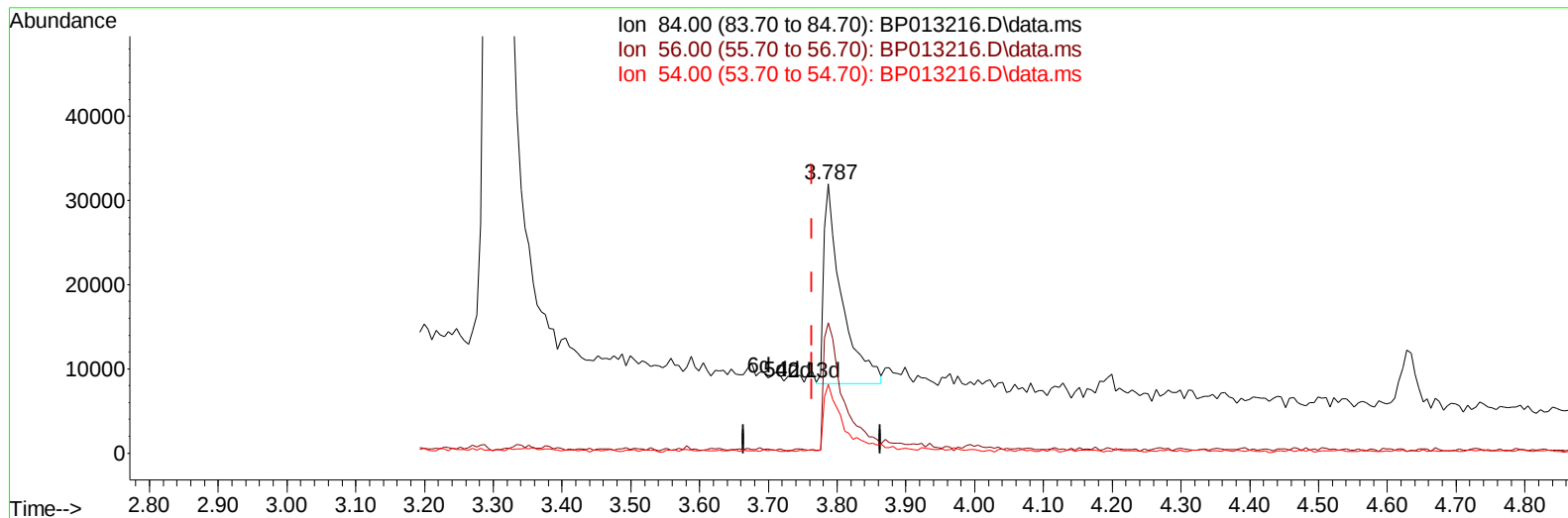
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TIC: BP013216.D\data.ms

(4) Pyridine-d5 (S)

3.787min (+ 0.024) 1.25 ng/ul m

response 42882

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	48.50#
54.00	30.50	25.61
0.00	0.00	0.00

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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	515066	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	2051682	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.593	164	1155464	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	2193409	20.000	ng/ul	-0.01
79) Chrysene-d12	21.433	240	2196978	20.000	ng/ul	0.00
88) Perylene-d12	23.910	264	2599594	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	1335778	110.886	ng/uL	-0.04
4) Pyridine-d5	3.787	84	42882m	1.253	ng/ul	0.02
7) Phenol-d5	7.111	99	231588	5.520	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	641559	25.350	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	575532	17.427	ng/ul	-0.01
15) 4-Methylphenol-d8	8.658	113	384562	11.187	ng/ul	-0.01
21) Nitrobenzene-d5	9.116	128	413331	27.420	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.840	143	418959	24.687	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.381	165	696774	21.138	ng/ul	0.00
31) 4-Chloroaniline-d4	10.905	131	411444	8.842	ng/ul	0.00
46) Dimethylphthalate-d6	13.999	166	2553947	28.760	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	2786442	28.556	ng/ul	0.00
54) 4-Nitrophenol-d4	14.793	143	51961	3.253	ng/ul	0.00
60) Fluorene-d10	15.587	176	2189045	29.138	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	281126	19.917	ng/ul	0.00
73) Anthracene-d10	17.445	188	3468595	35.052	ng/ul	-0.01
81) Pyrene-d10	19.681	212	4167619	33.048	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	4534052	34.992	ng/ul	-0.01

Target Compounds Qvalue

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

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