

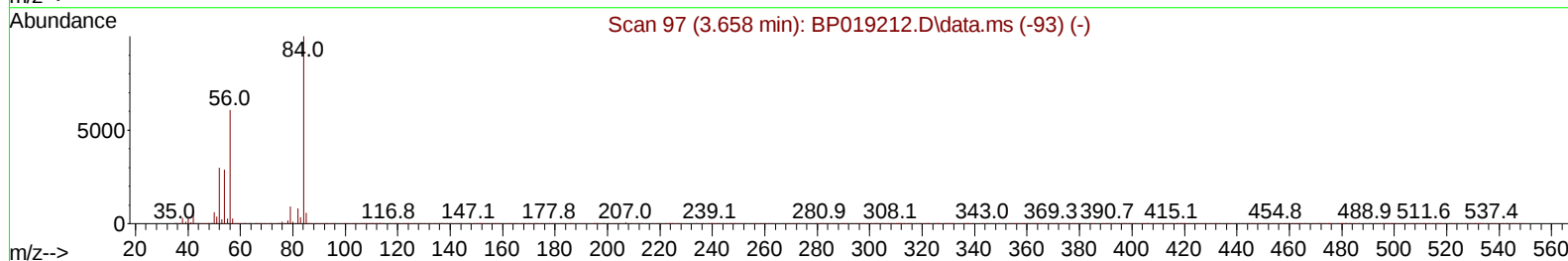
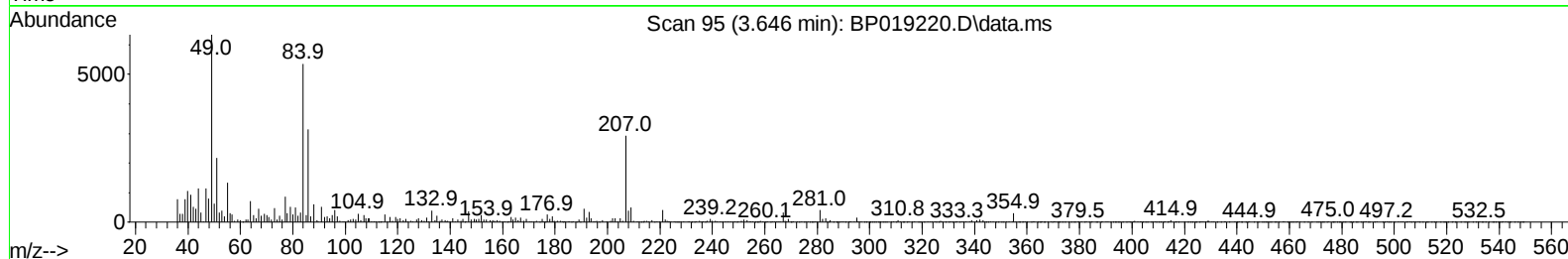
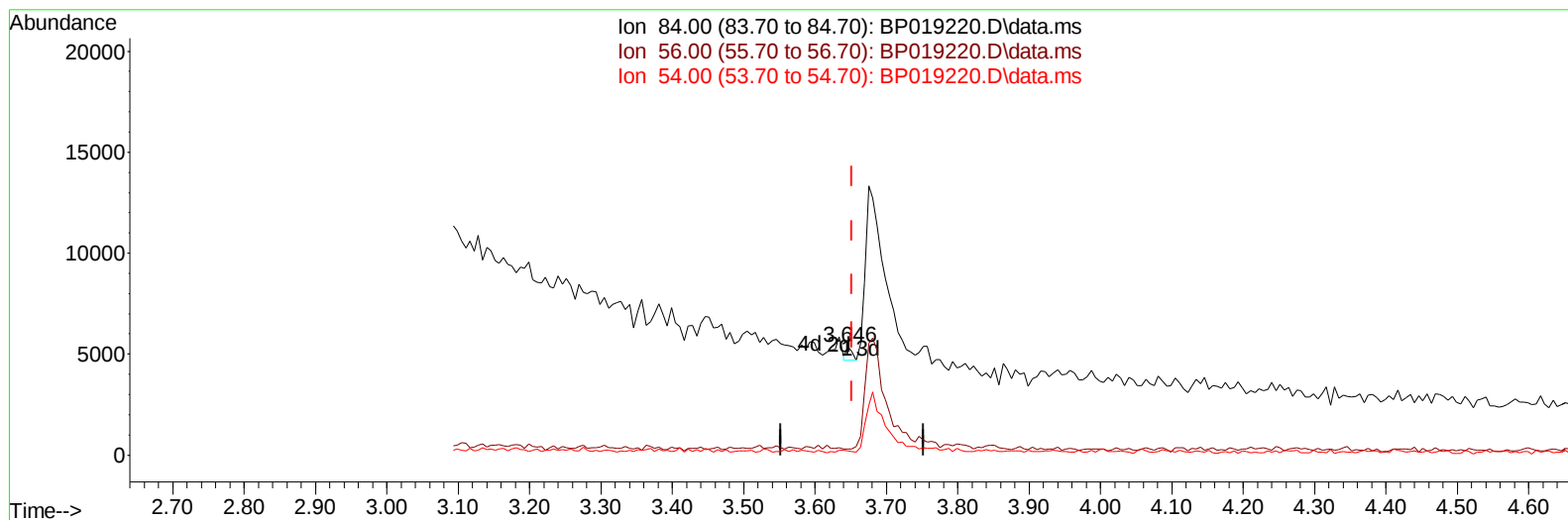
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010224\
 Data File : BP019220.D
 Acq On : 02 Jan 2024 22:01
 Operator : MA/JU
 Sample : 06020-03RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH3P1RE

Manual IntegrationsAPPROVED

Quant Time: Jan 03 09:35:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/04/2024
 Supervised By :mohammad ahmed 01/05/2024



TIC: BP019220.D\data.ms

(4) Pyridine-d5 (S)

3.646min (-0.006) 0.02 ng/u1

response 376

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	5.72#
54.00	29.60	3.58#
0.00	0.00	0.00

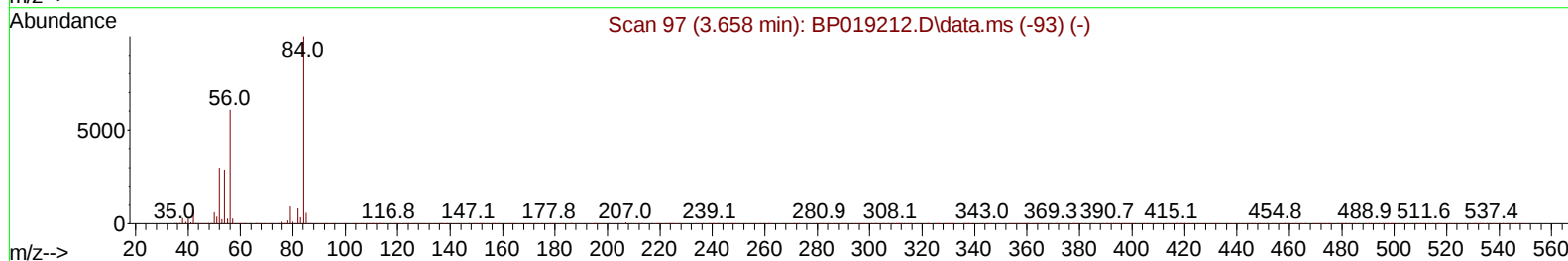
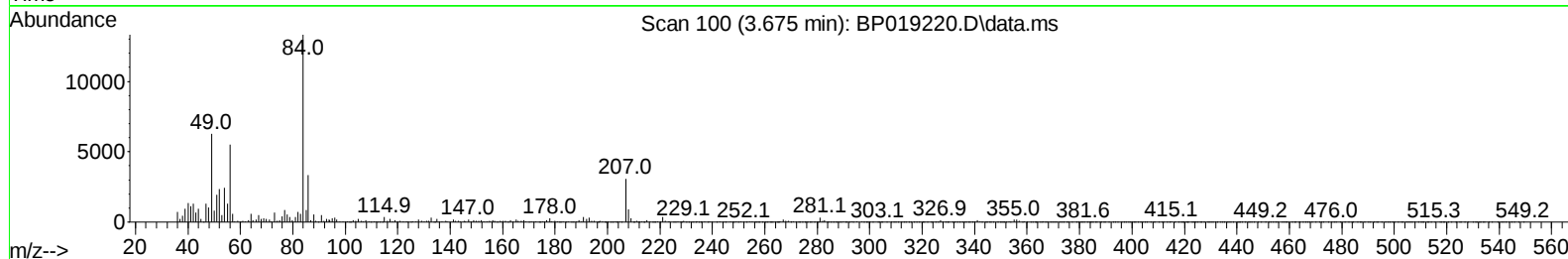
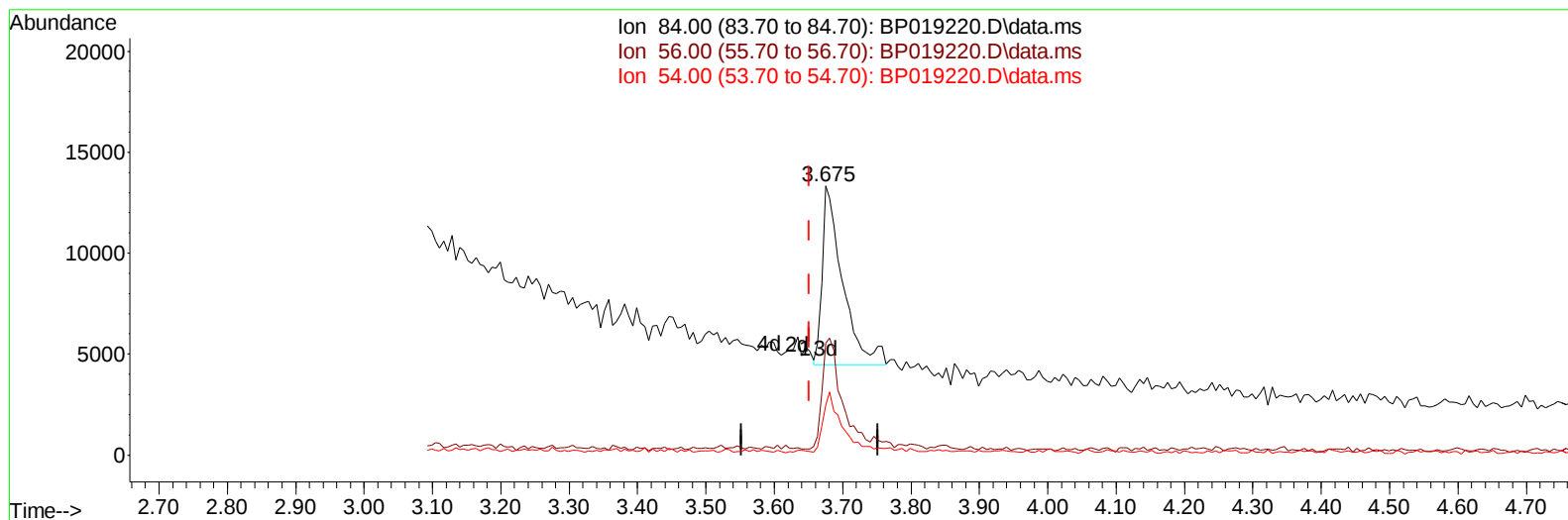
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 Operator : MA/JU
 Sample : 06020-03RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

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 ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Jan 02 23:04:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
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TIC: BP019220.D\data.ms

(4) Pyridine-d5 (S)

3.675min (+ 0.023) 1.09 ng/u1 m

response 18346

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.10	41.53#
54.00	29.60	18.41#
0.00	0.00	0.00

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 Operator : MA/JU
 Sample : 06020-03RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	229539	20.000	ng/ul	0.00
20) Naphthalene-d8	10.592	136	904815	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	577402	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	1374074	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1428596	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1474055	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	30035	4.469	ng/uL	0.00
4) Pyridine-d5	3.675	84	18346m	1.088	ng/ul	0.02
7) Phenol-d5	6.981	99	158526	7.701	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	488702	41.038	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	544222	32.795	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	231362	14.437	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	327385	41.856	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	371211	43.914	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	656829	38.480	ng/ul	0.00
31) 4-Chloroaniline-d4	10.751	131	16730	0.797	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1917905	37.605	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	205388	3.711	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	102875	11.945	ng/ul	0.01
60) Fluorene-d10	15.439	176	1546186	35.093	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	351927	39.068	ng/ul	0.00
73) Anthracene-d10	17.298	188	1314149	18.314	ng/ul	0.00
81) Pyrene-d10	19.545	212	2356932	27.071	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	848767	10.017	ng/ul	0.00

Target Compounds Qvalue

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

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