

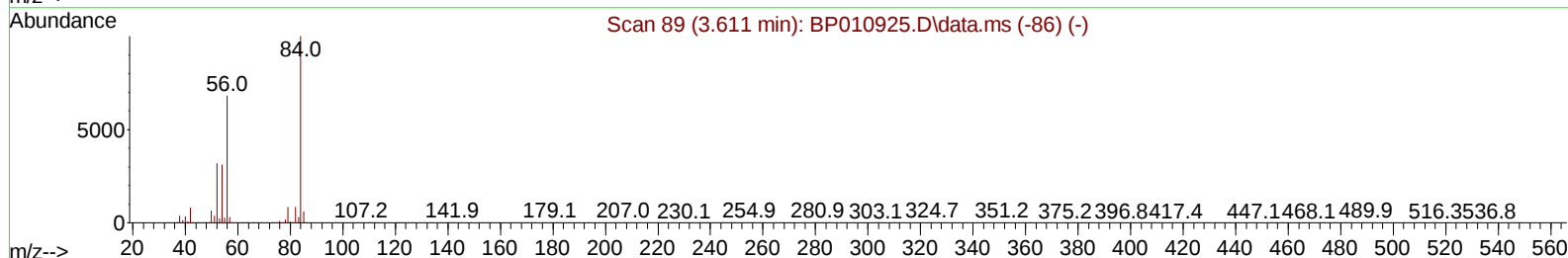
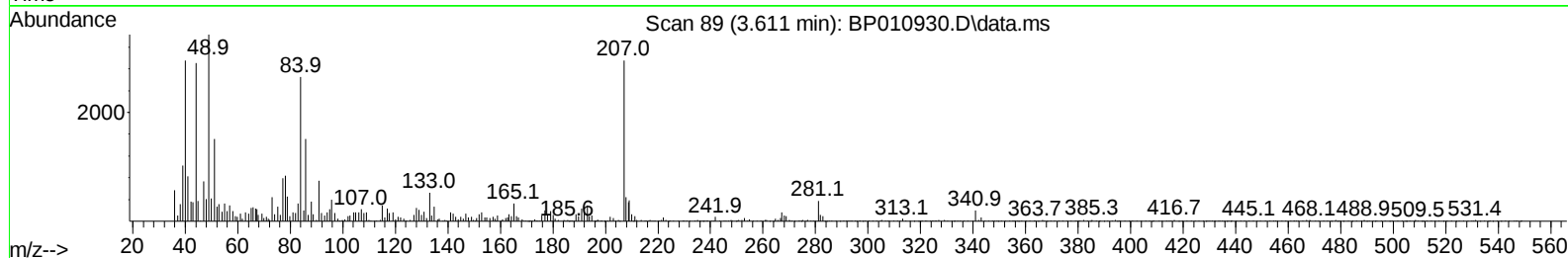
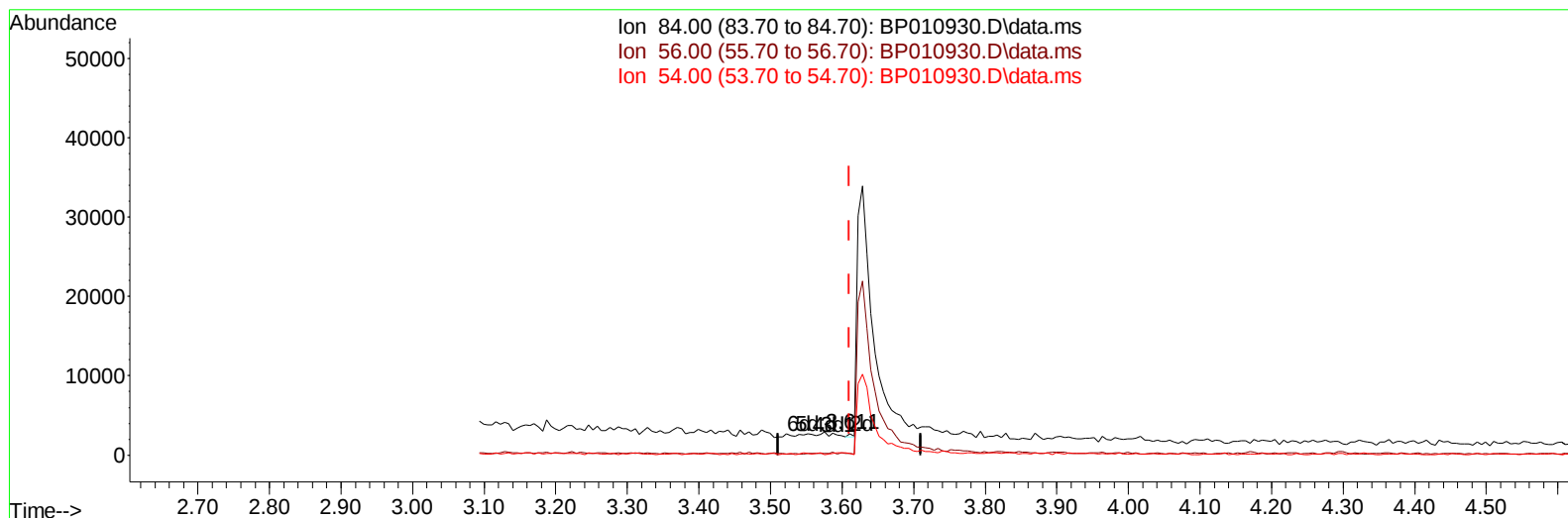
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070622\
 Data File : BP010930.D
 Acq On : 06 Jul 2022 15:18
 Operator : CG/JU
 Sample : N3555-01DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA8DL

Manual IntegrationsAPPROVED

Quant Time: Jul 06 23:15:57 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 06 23:13:53 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/07/2022
 Supervised By :mohammad ahmed 07/07/2022



TIC: BP010930.D\data.ms

(4) Pyridine-d5 (S)

3.611min (+ 0.000) 0.01 ng/u1

response 151

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	7.05#
54.00	44.80	6.41#
0.00	0.00	0.00

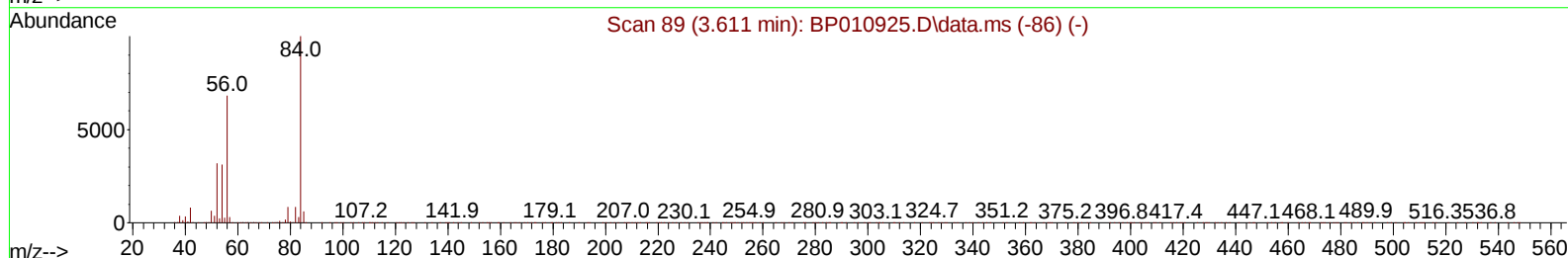
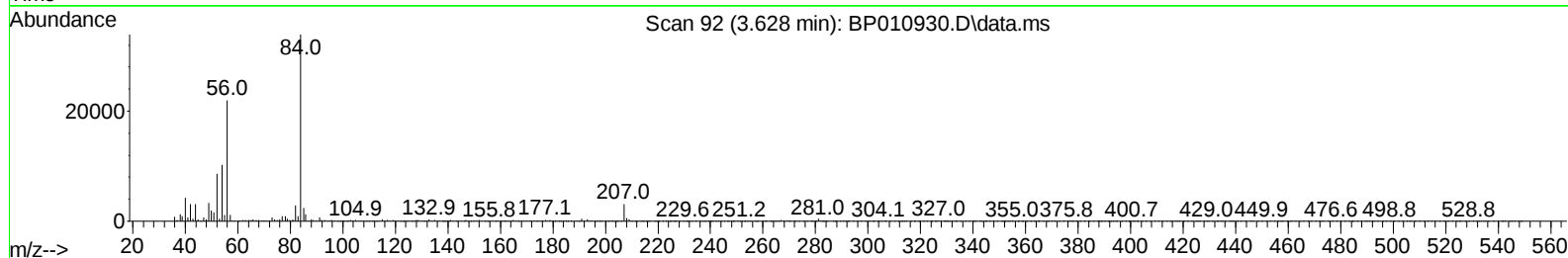
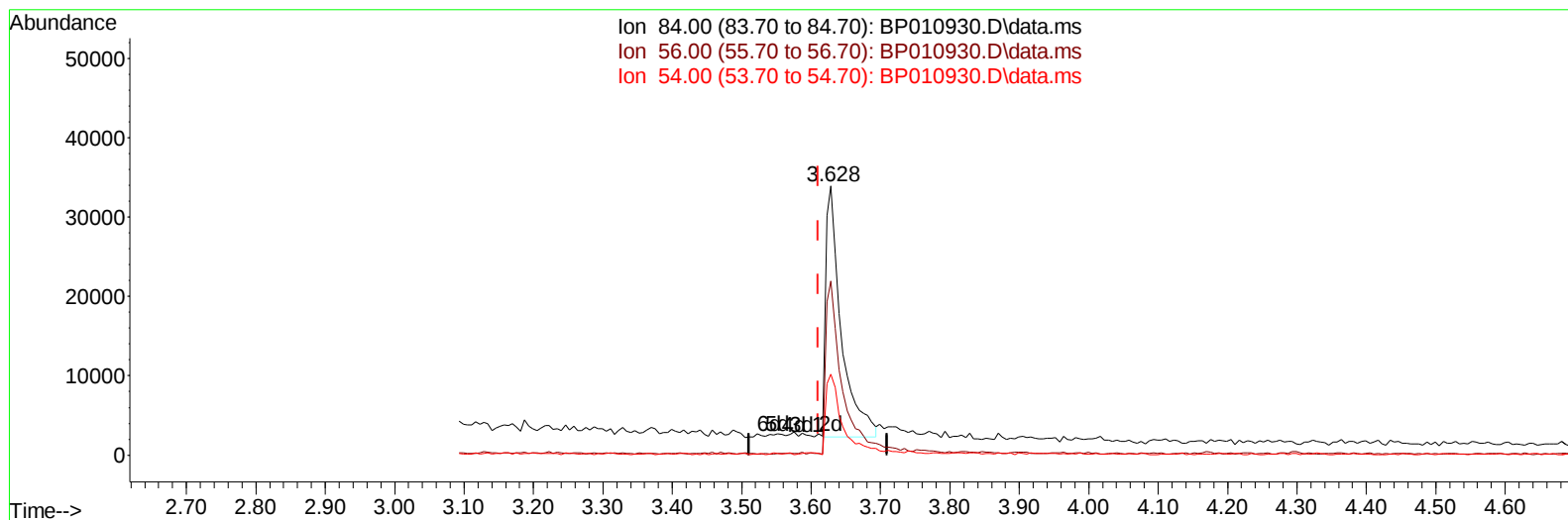
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TIC: BP010930.D\data.ms

(4) Pyridine-d5 (S)

3.628min (+ 0.018) 1.74 ng/ul m

response 48705

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	64.81#
54.00	44.80	30.05#
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	7.711	152	398574	20.000	ng/ul	0.00
20) Naphthalene-d8	10.505	136	1511089	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	744010	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.134	188	1402781	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.245	240	1424119	20.000	ng/ul	# 0.00
88) Perylene-d12	23.610	264	1579838	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	11598	1.101	ng/uL	0.00
4) Pyridine-d5	3.628	84	48705m	1.737	ng/ul	0.02
7) Phenol-d5	6.887	99	44945	1.425	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	126032	6.108	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	124187	4.744	ng/ul	0.00
15) 4-Methylphenol-d8	8.422	113	76184	3.054	ng/ul	0.00
21) Nitrobenzene-d5	8.869	128	67300	6.996	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	53202	6.685	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	104741	5.326	ng/ul	0.00
31) 4-Chloroaniline-d4	10.652	131	63394	1.975	ng/ul	0.01
46) Dimethylphthalate-d6	13.787	166	338383	6.640	ng/ul	0.00
49) Acenaphthylene-d8	14.057	160	390424	6.002	ng/ul	0.00
54) 4-Nitrophenol-d4	14.581	143	8535	1.072	ng/ul	0.00
60) Fluorene-d10	15.363	176	305254	6.979	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.487	200	20272	4.362	ng/ul	0.00
73) Anthracene-d10	17.228	188	443253	7.102	ng/ul	0.00
81) Pyrene-d10	19.481	212	498287	6.422	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.451	264	544952	6.980	ng/ul	0.00
Target Compounds						
					Qvalue	
52) Acenaphthene	14.428	153	1903384	41.409	ng/ul	99
61) Fluorene	15.422	166	651382	12.975	ng/ul	99
74) Anthracene	17.263	178	117818	1.568	ng/ul	96
77) Carbazole	17.545	167	293577	4.341	ng/ul	100
80) Fluoranthene	19.157	202	289853	3.038	ng/ul#	91
82) Pyrene	19.510	202	168487	1.748	ng/ul#	86

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

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