

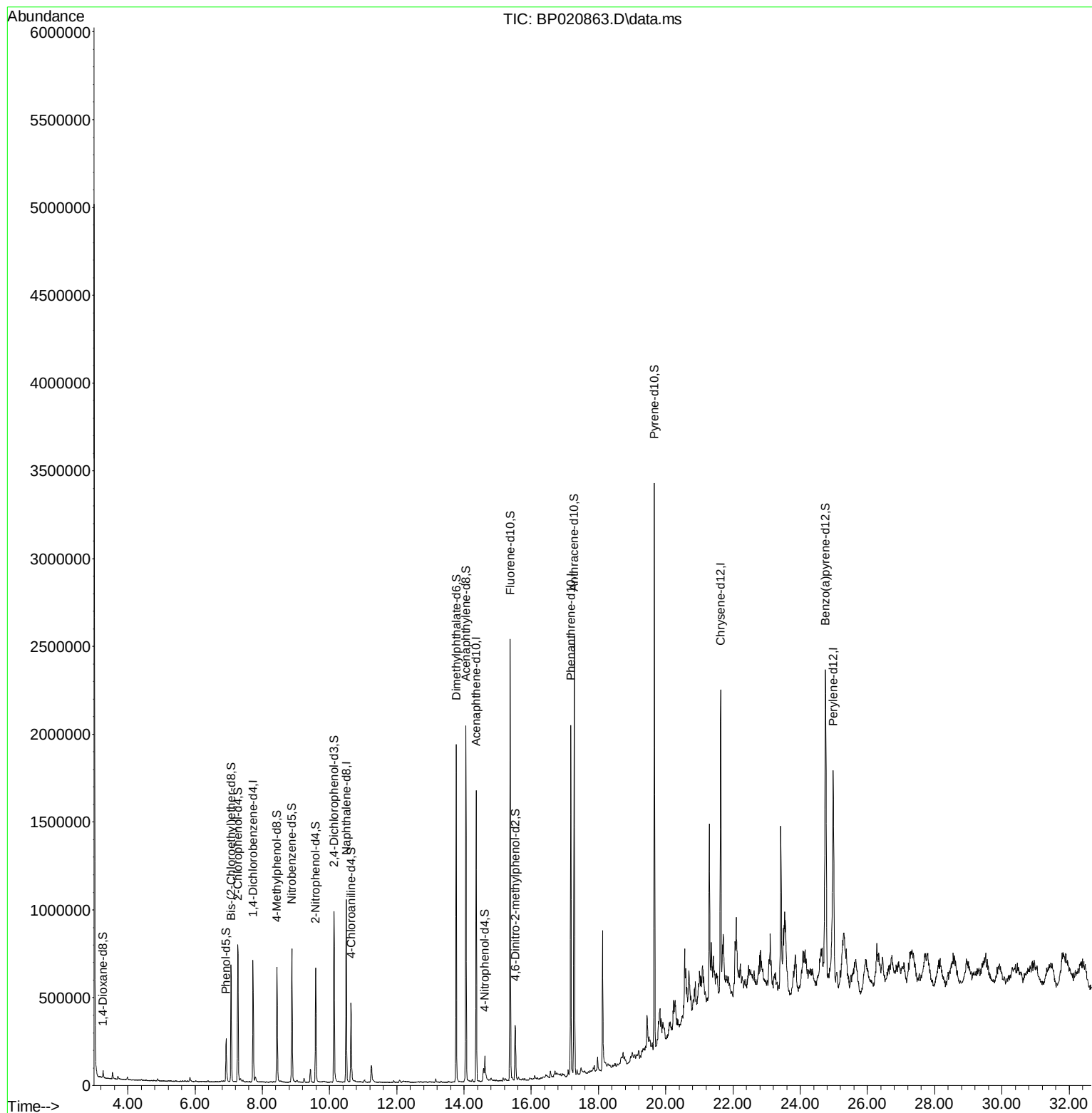
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020863.D
Acq On : 09 Jul 2024 14:01
Operator : MA/JU
Sample : P3084-07
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1G89

Manual IntegrationsAPPROVED

Quant Time: Jul 09 23:06:43 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 03 15:01:56 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/10/2024
Supervised By :mohammad ahmed 07/10/2024



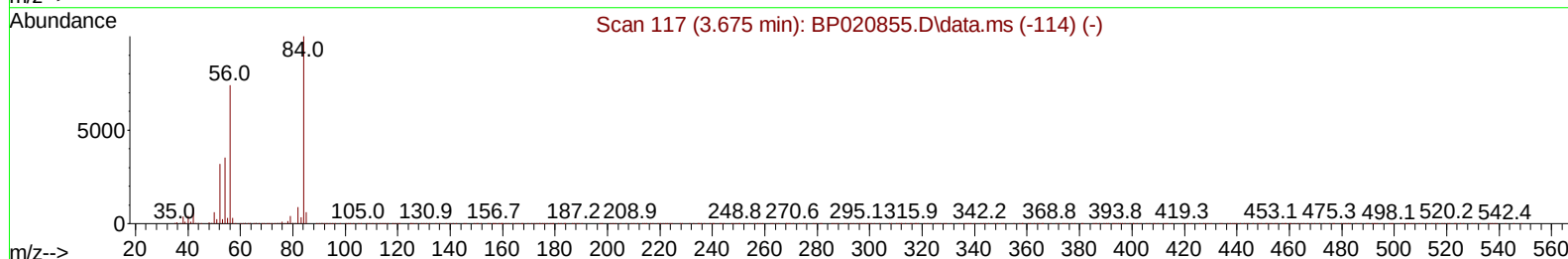
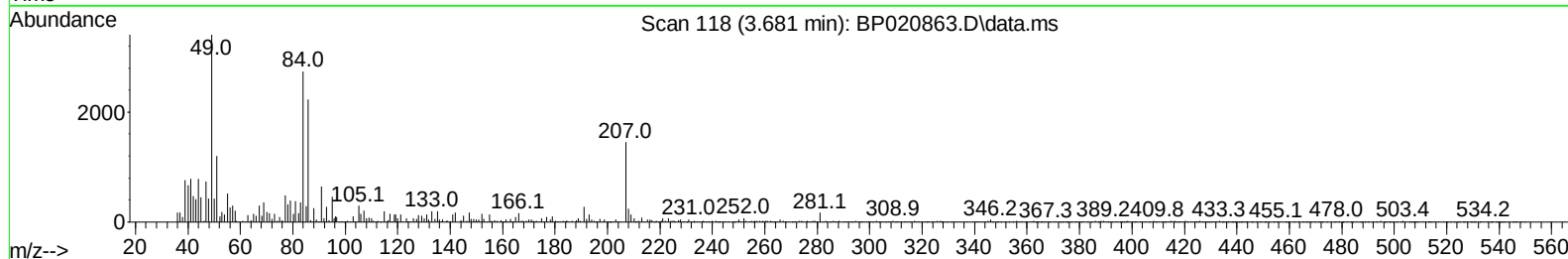
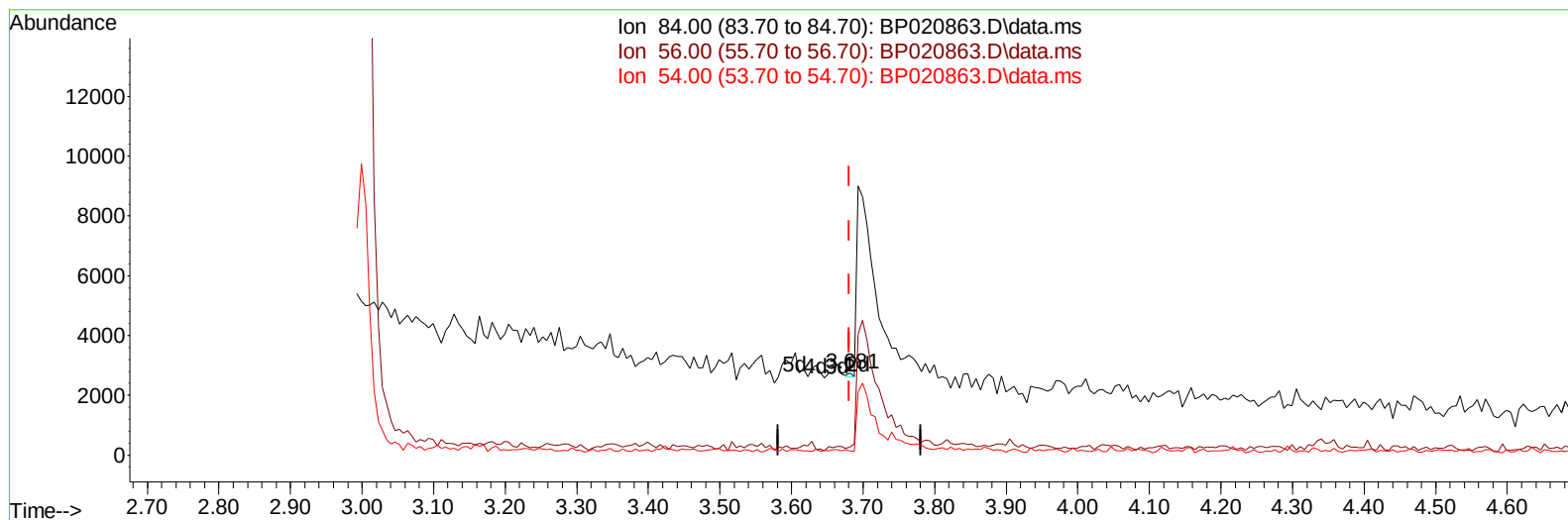
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TIC: BP020863.D\data.ms

(4) Pyridine-d5 (S)

3.681min (+ 0.000) 0.00 ng/u1

response	39	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	75.20	9.56
54.00	37.40	4.96
0.00	0.00	0.00

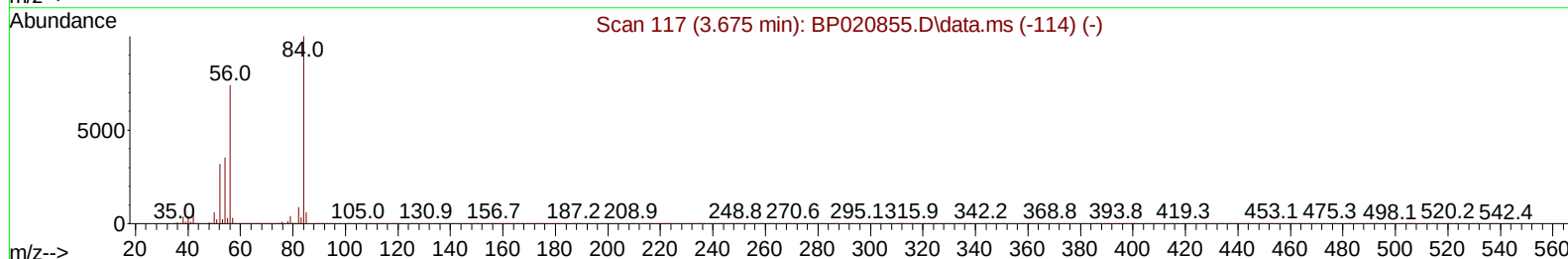
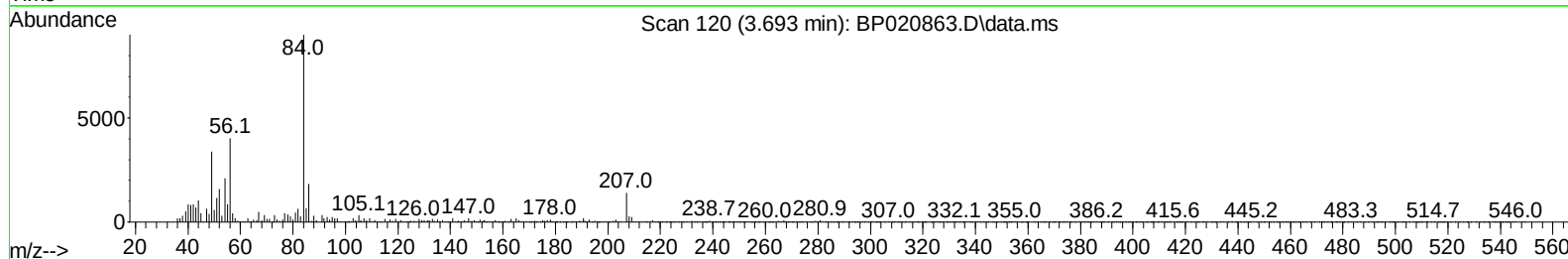
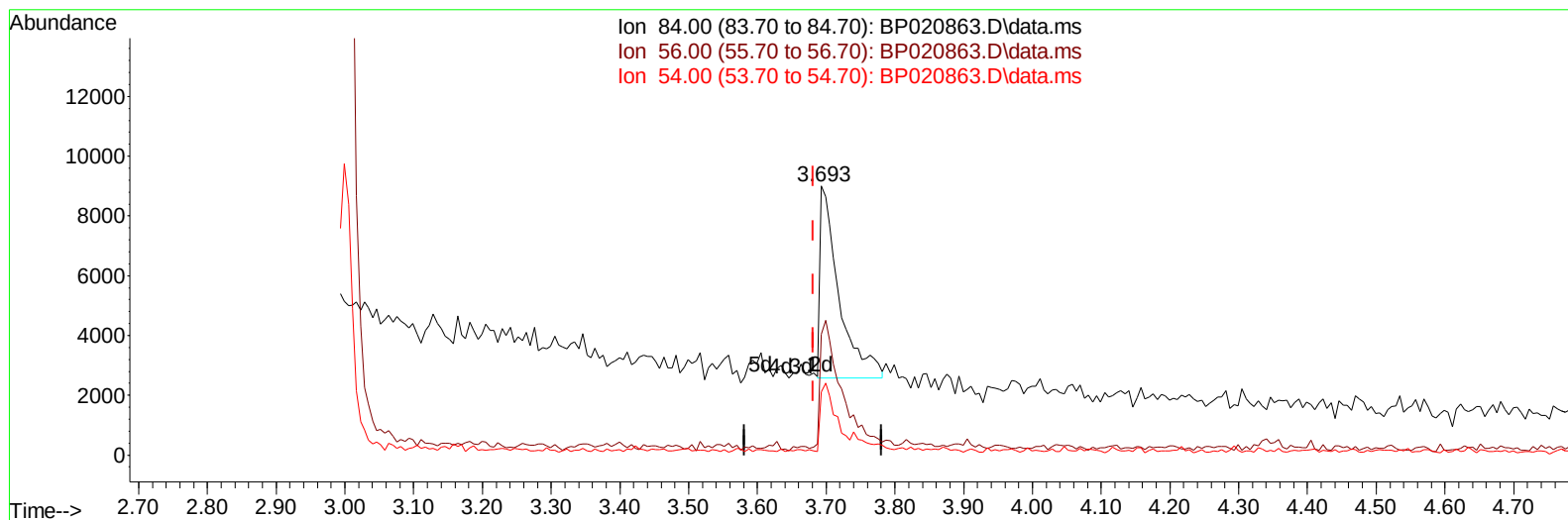
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(4) Pyridine-d5 (S)

3.693min (+ 0.012) 0.92 ng/ul m

response 12392

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	75.20	44.89#
54.00	37.40	23.37#
0.00	0.00	0.00

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 ALS Vial : 37 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	198379	20.000	ng/ul	0.00
20) Naphthalene-d8	10.499	136	817638	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.363	164	551736	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.175	188	1168713	20.000	ng/ul	0.01
79) Chrysene-d12	21.633	240	1056339	20.000	ng/ul	0.01
88) Perylene-d12	24.980	264	1238042	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	20497	4.437	ng/uL	0.00
4) Pyridine-d5	3.693	84	12392m	0.920	ng/ul	0.01
7) Phenol-d5	6.922	99	152396	9.366	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.069	67	280368	27.181	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.269	132	363525	27.337	ng/ul	0.00
15) 4-Methylphenol-d8	8.434	113	268157	20.292	ng/ul	0.00
21) Nitrobenzene-d5	8.881	128	192455	31.881	ng/ul	0.00
24) 2-Nitrophenol-d4	9.587	143	222327	36.803	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.134	165	377375	30.327	ng/ul	0.00
31) 4-Chloroaniline-d4	10.640	131	279015	16.000	ng/ul	0.00
46) Dimethylphthalate-d6	13.769	166	1179751	30.753	ng/ul	0.01
49) Acenaphthylene-d8	14.057	160	1339672	29.266	ng/ul	0.02
54) 4-Nitrophenol-d4	14.616	143	49779	8.493	ng/ul	0.04
60) Fluorene-d10	15.369	176	1010871	31.315	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.522	200	87773	14.903	ng/ul	0.02
73) Anthracene-d10	17.275	188	1664909	31.742	ng/ul	0.01
81) Pyrene-d10	19.663	212	1932217	30.448	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.751	264	1907085	31.667	ng/ul	0.01

Target Compounds Qvalue

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