

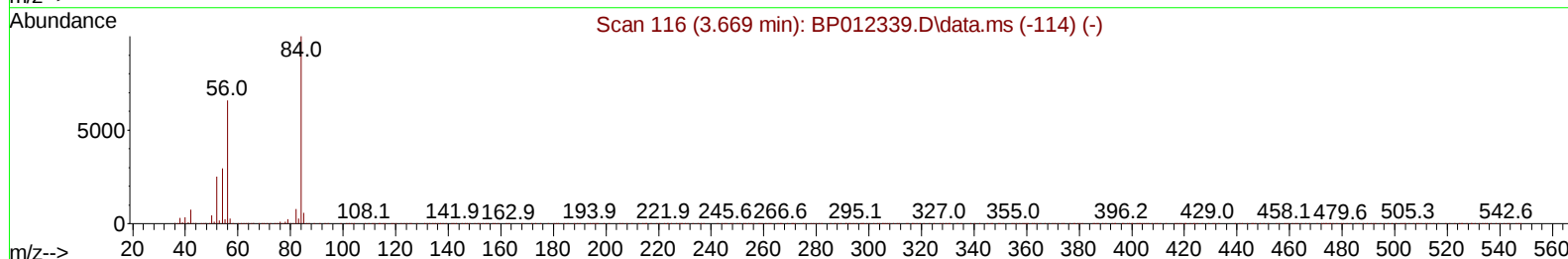
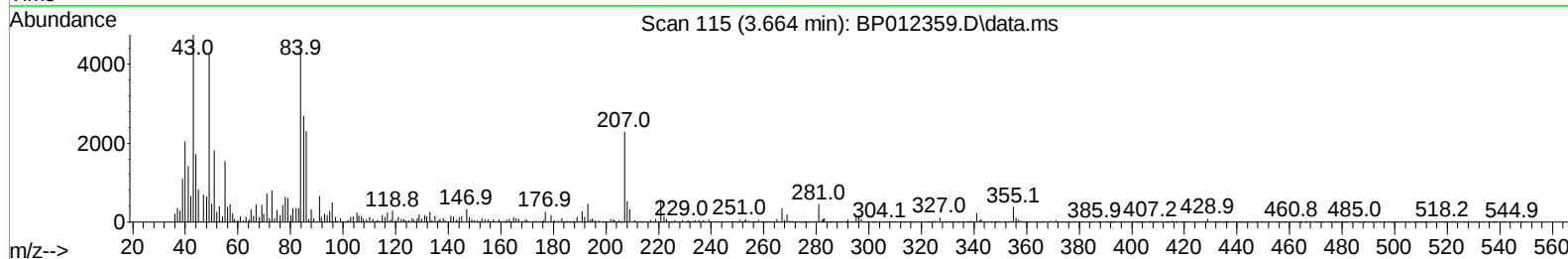
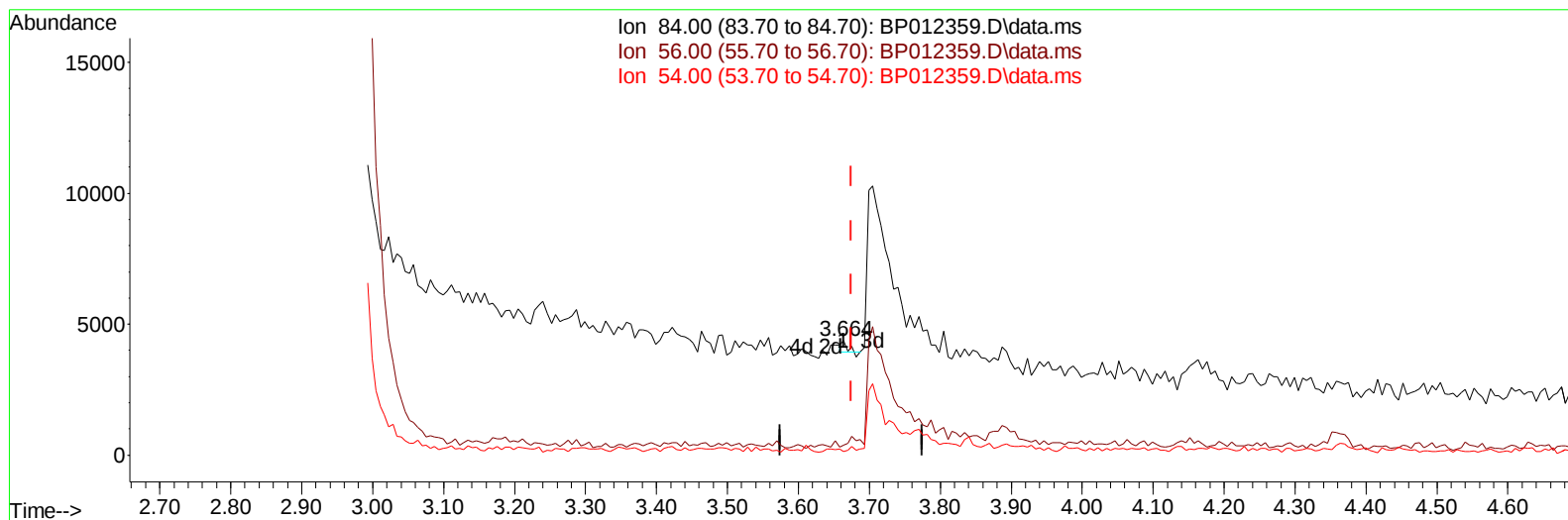
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012359.D
 Acq On : 28 Oct 2022 00:40
 Operator : CG/JU
 Sample : N5144-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBLO

Manual IntegrationsAPPROVED

Quant Time: Oct 28 01:39:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 26 23:40:20 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
 Supervised By :mohammad ahmed 10/31/2022



TIC: BP012359.D\data.ms

(4) Pyridine-d5 (S)

3.664min (-0.011) 0.01 ng/u1

response	171	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	8.94#
54.00	29.80	3.28#
0.00	0.00	0.00

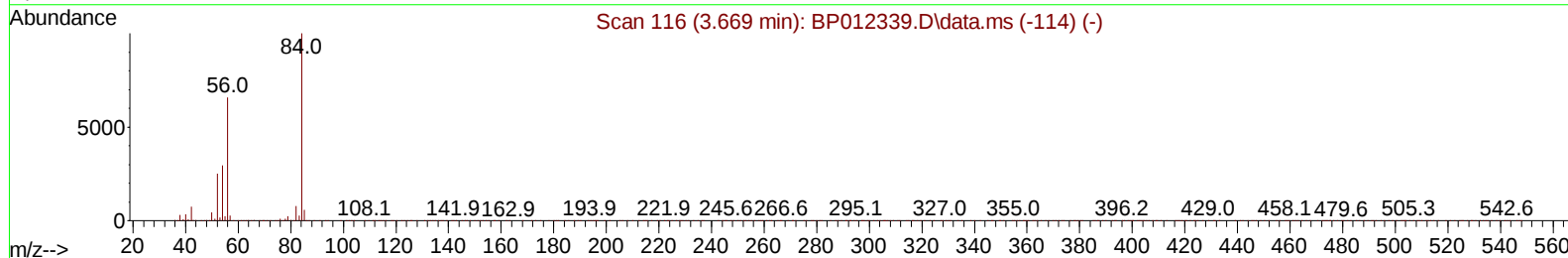
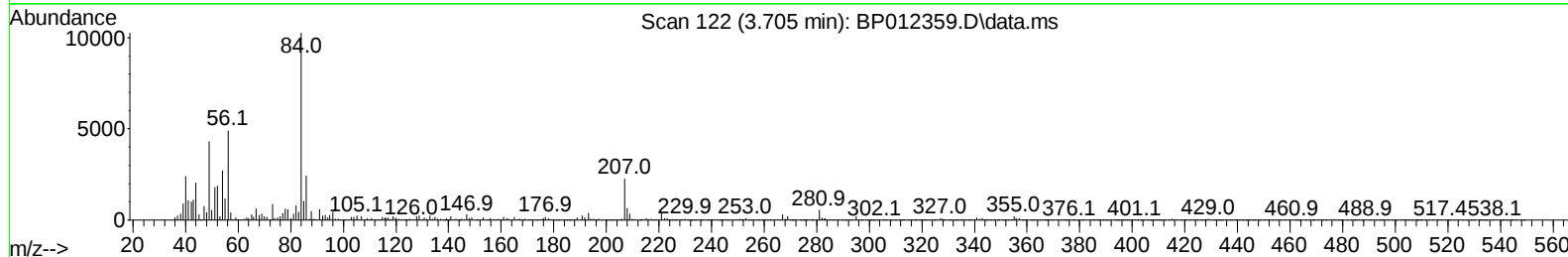
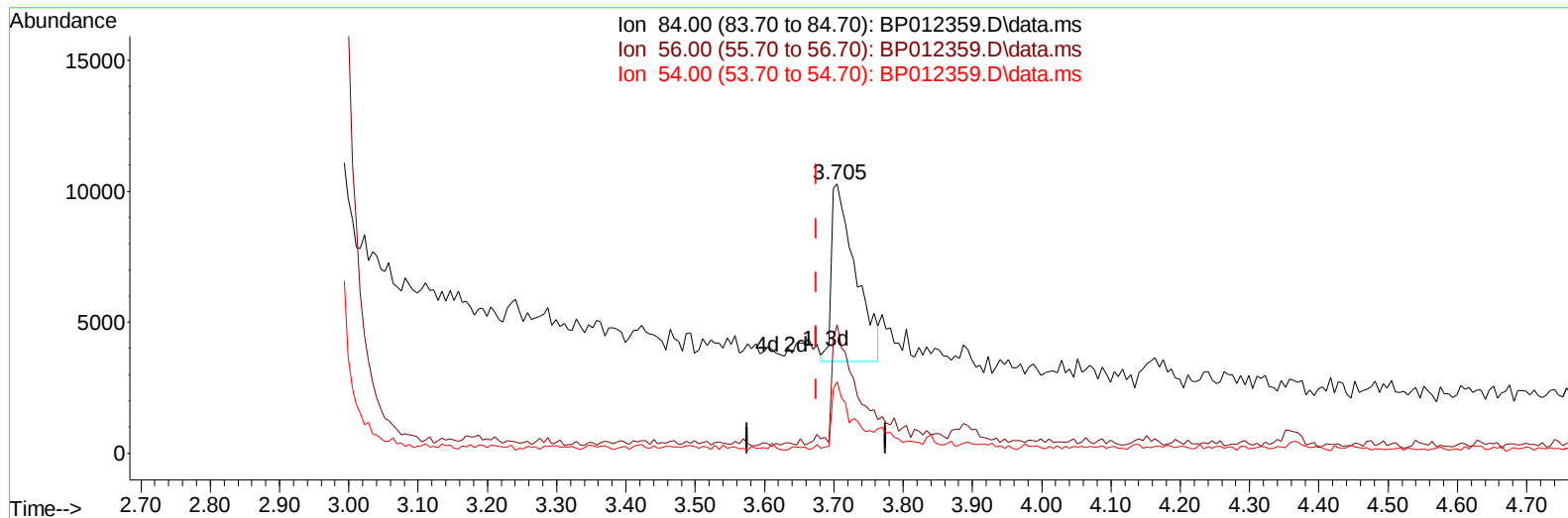
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(4) Pyridine-d5 (S)

3.705min (+ 0.030) 0.68 ng/ul m

response 16344

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	47.88#
54.00	29.80	26.51
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	7.805	152	337935	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.610	136	1583391	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	1077106	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	2347112	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	2437870	20.000	ng/ul	0.00
88) Perylene-d12	23.710	264	2268360	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	33360	3.975	ng/uL	0.00
4) Pyridine-d5	3.705	84	16344m	0.678	ng/ul	0.03
7) Phenol-d5	6.975	99	680973	22.467	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	435887	24.090	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	552348	24.954	ng/ul	-0.01
15) 4-Methylphenol-d8	8.522	113	334000	13.933	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	302232	29.434	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	335699	31.499	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	567760	24.582	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	685082	19.921	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	2038400	27.645	ng/ul	-0.01
49) Acenaphthylene-d8	14.151	160	2301457	27.313	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	311629	22.331	ng/ul	0.00
60) Fluorene-d10	15.451	176	1796791	28.462	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.587	200	245353	20.078	ng/ul	0.00
73) Anthracene-d10	17.316	188	2715383	26.850	ng/ul	0.00
81) Pyrene-d10	19.557	212	3481121	28.441	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3114826	28.103	ng/ul	0.00

Target Compounds Qvalue

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

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