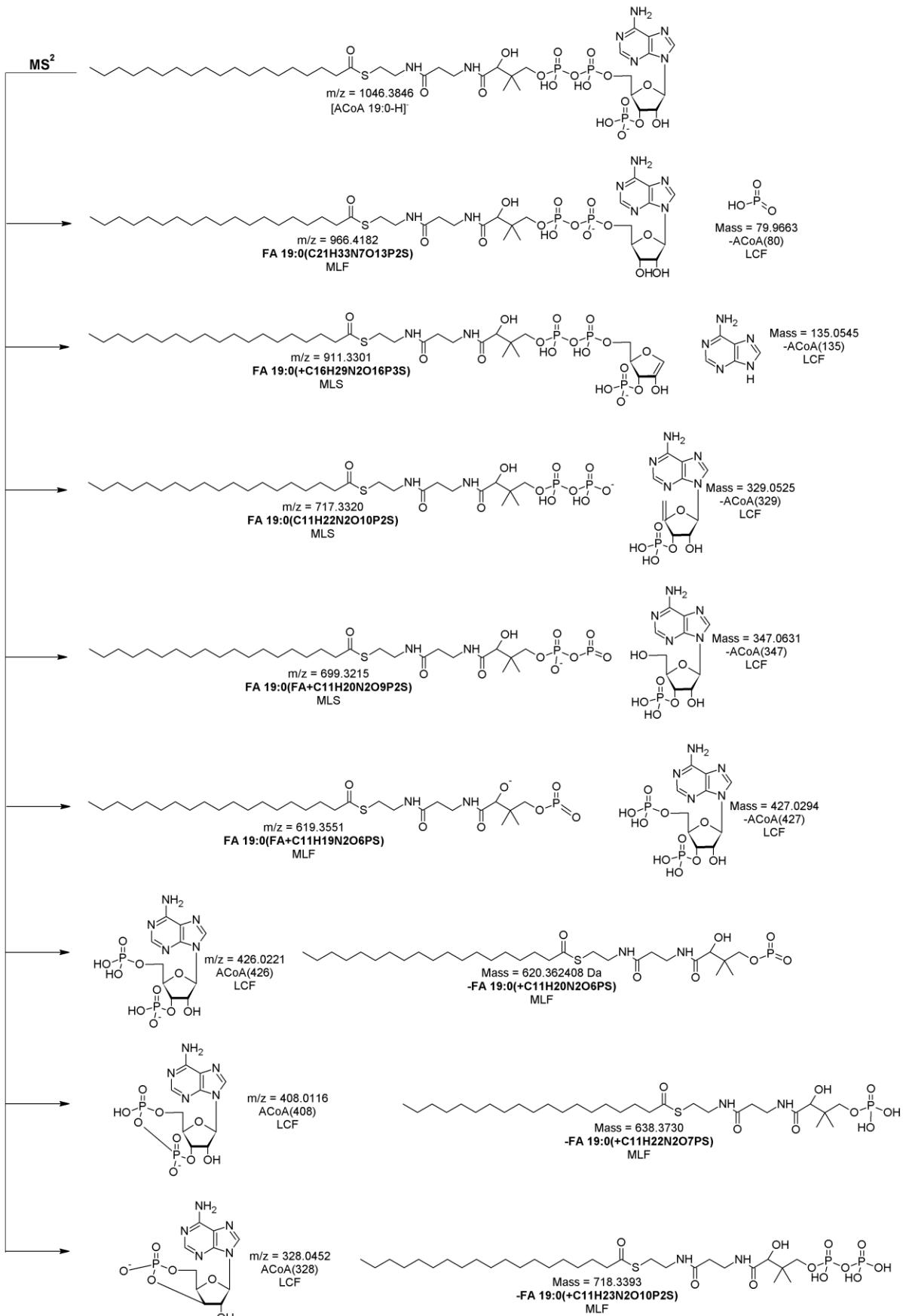
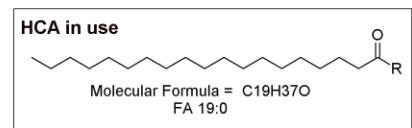


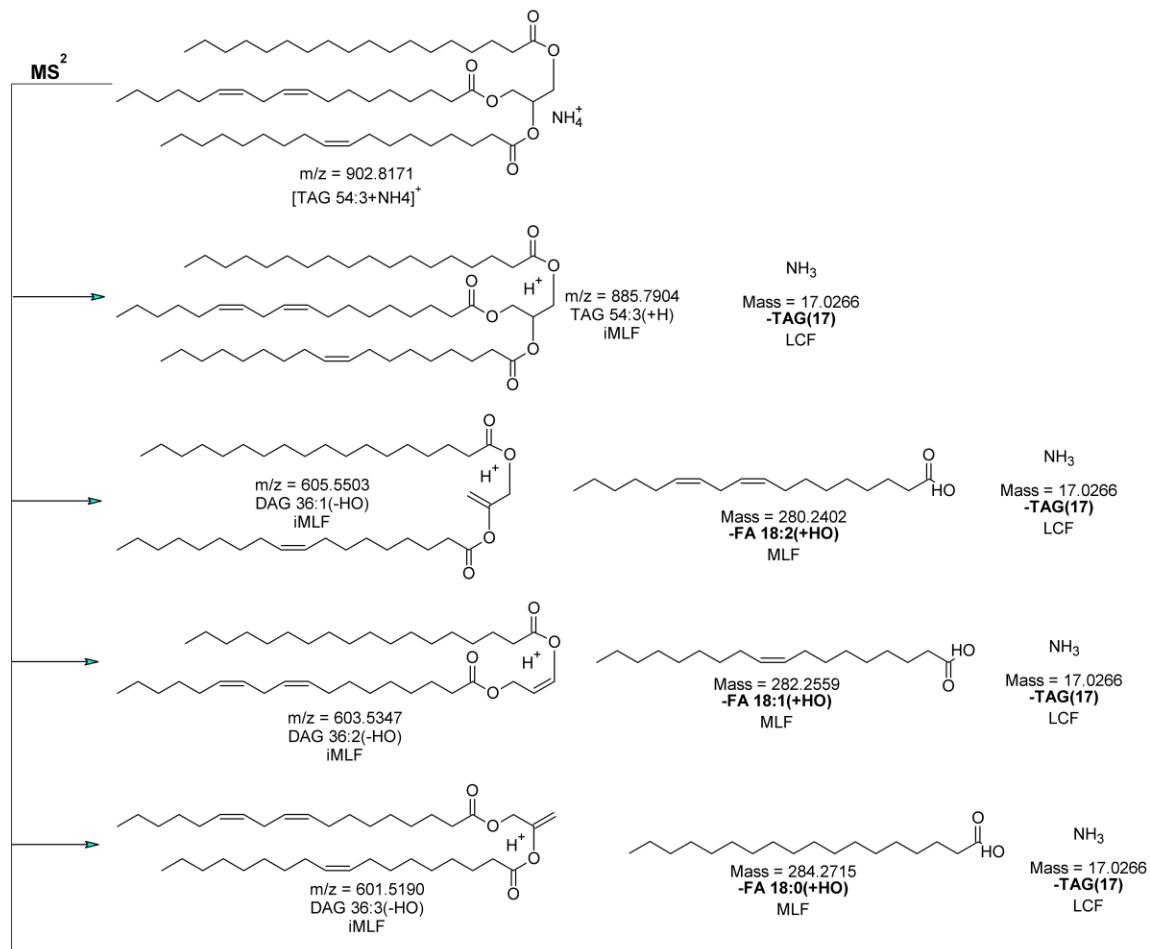
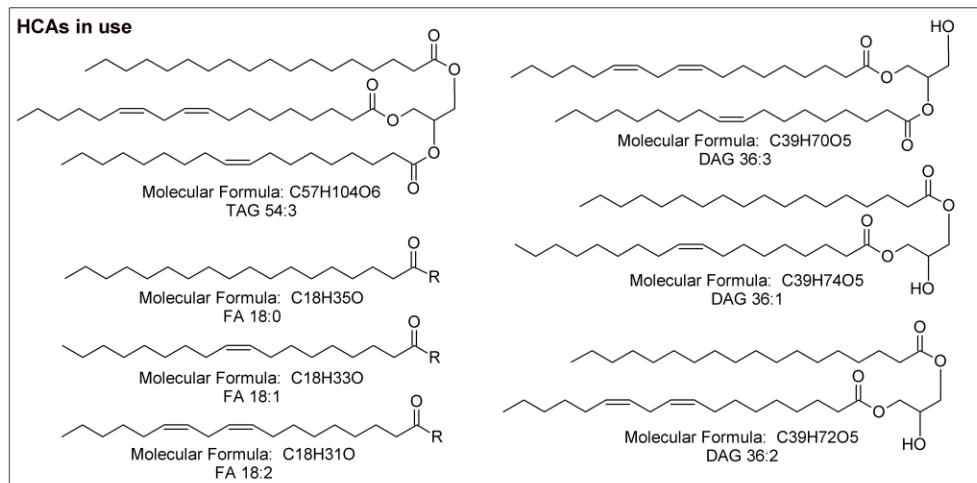
# S3A Fig) ACoA 19:0

Proposed structures corresponding to spectrum shown in Fig. 3A.

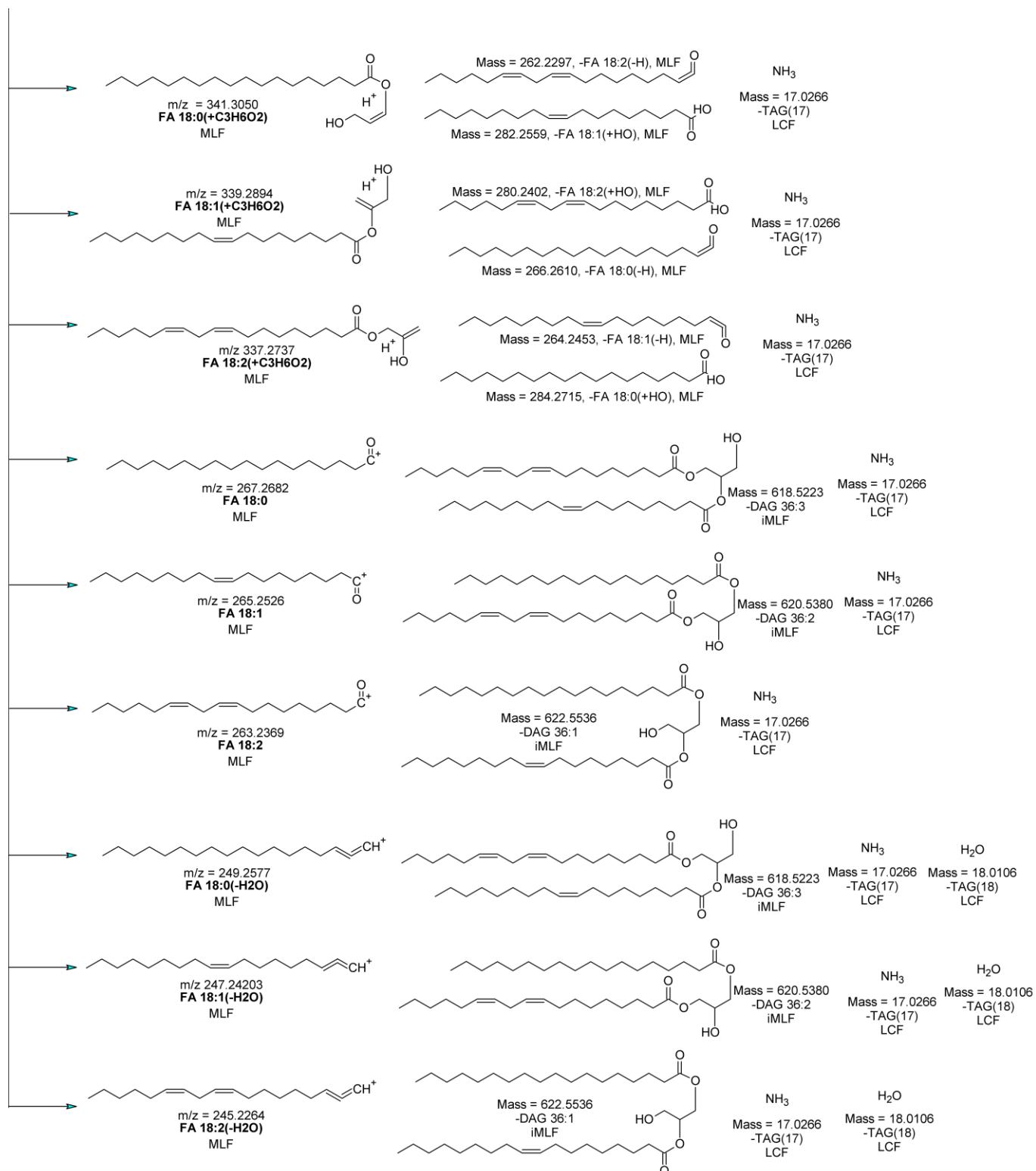


# S3B Fig) TAG 18:0-18:1-18:2

Proposed structures corresponding to spectrum shown in Fig. 3B.

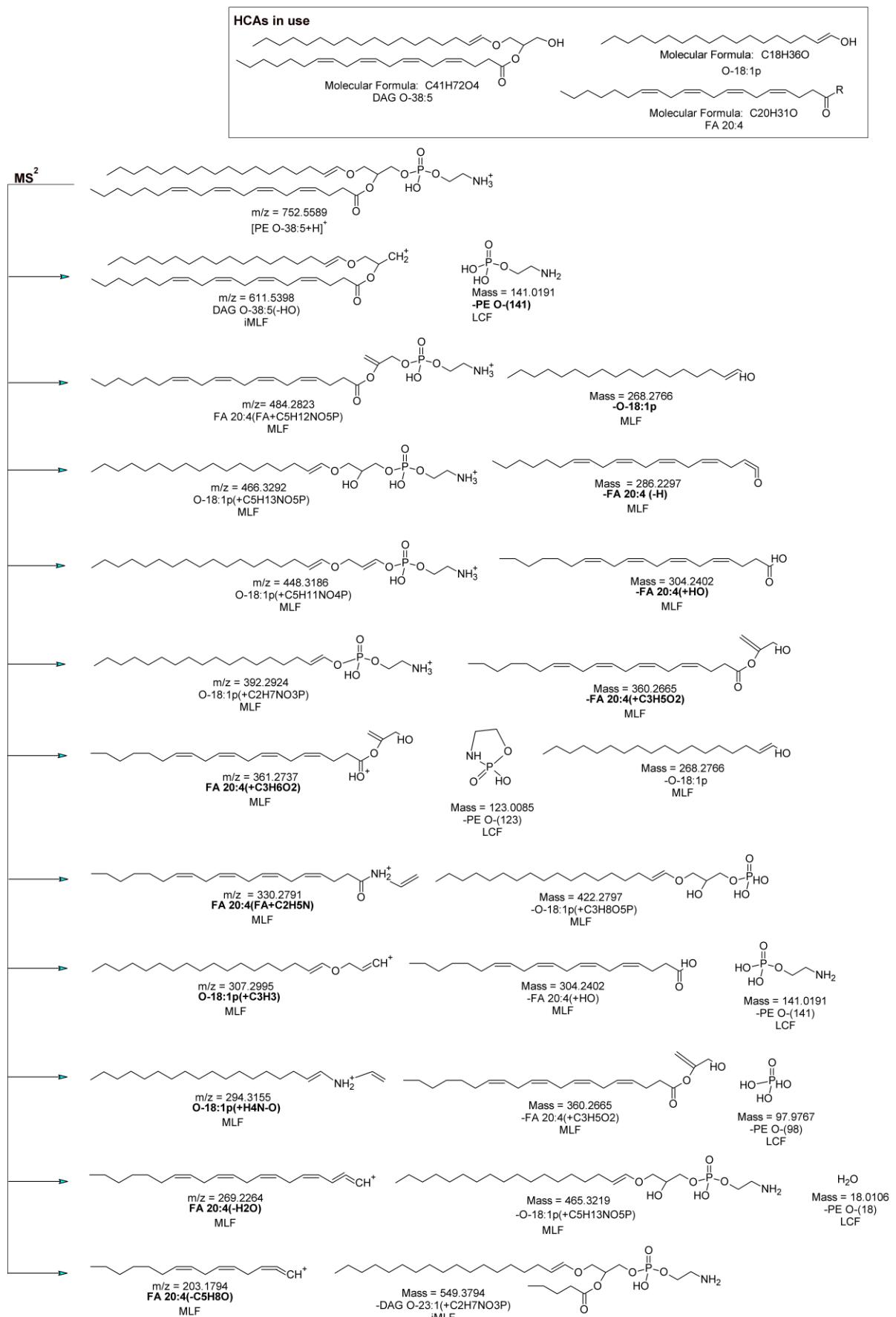


# S3B Fig) TAG 18:0-18:1-18:2 (cont.)



# S3C Fig) PE O-18:1p/20:4

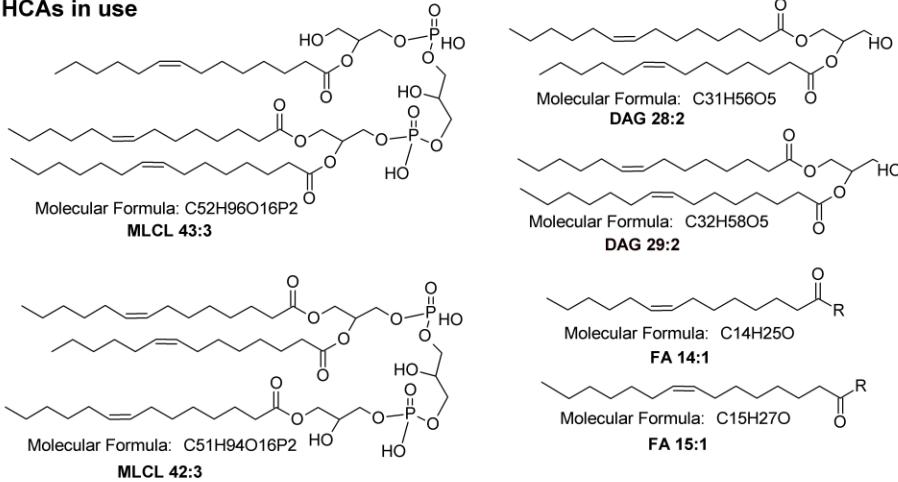
Proposed structures corresponding to spectrum shown in Fig. 3C.



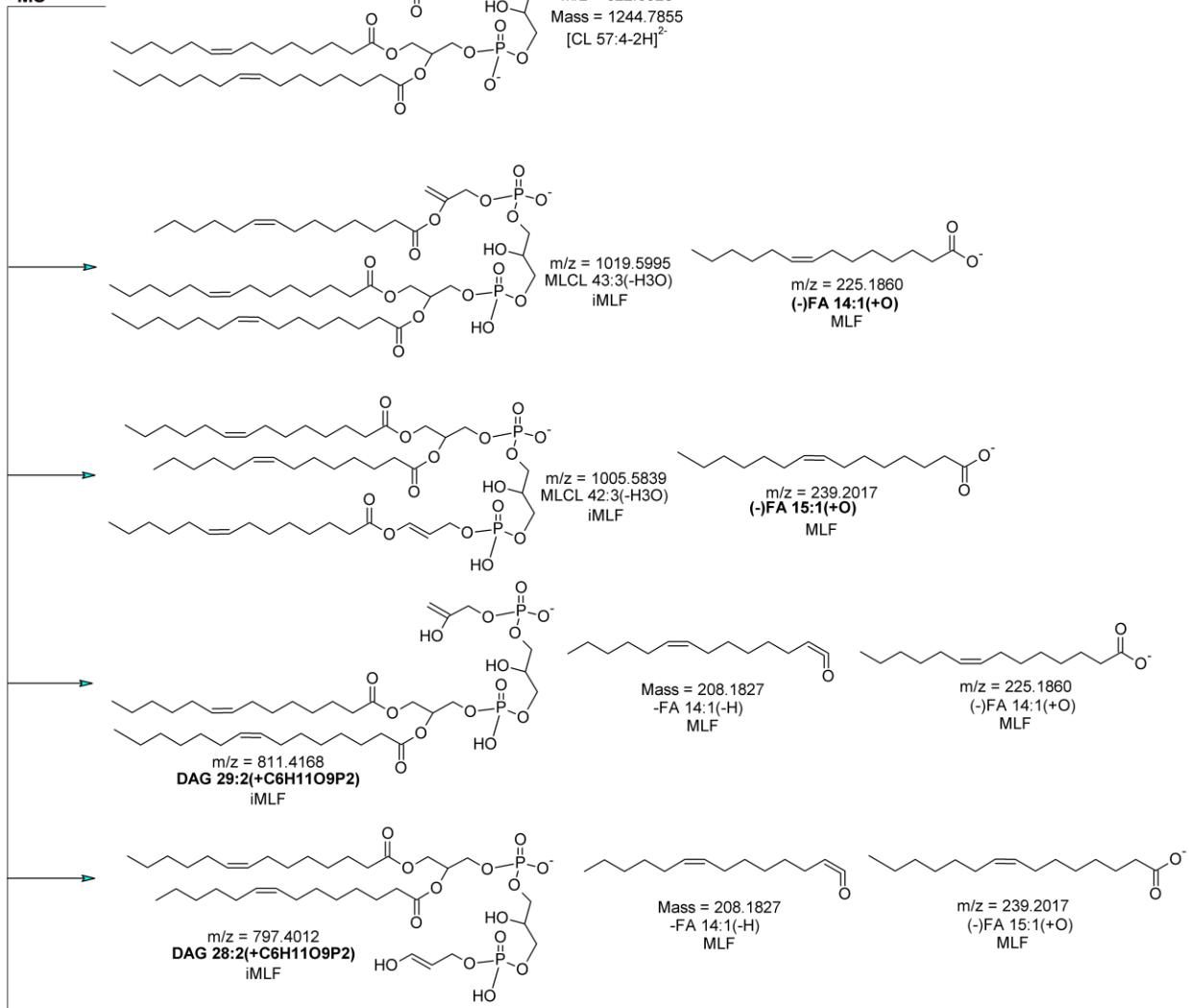
# S3D Fig) CL 14:1-14:1-14:1-15:1

Proposed structures corresponding to spectrum shown in Fig. 3D.

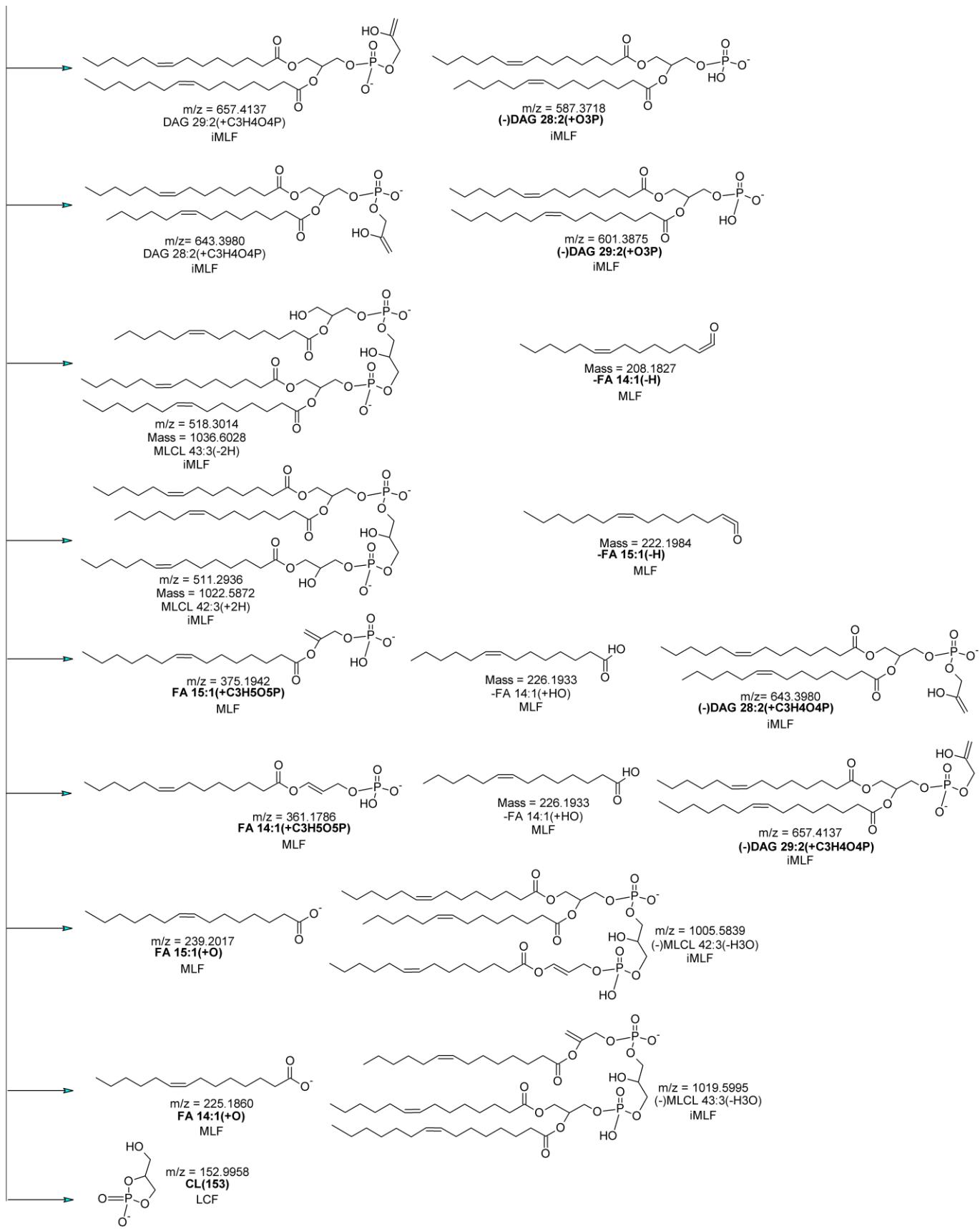
## HCAs in use



## MS<sup>2</sup>

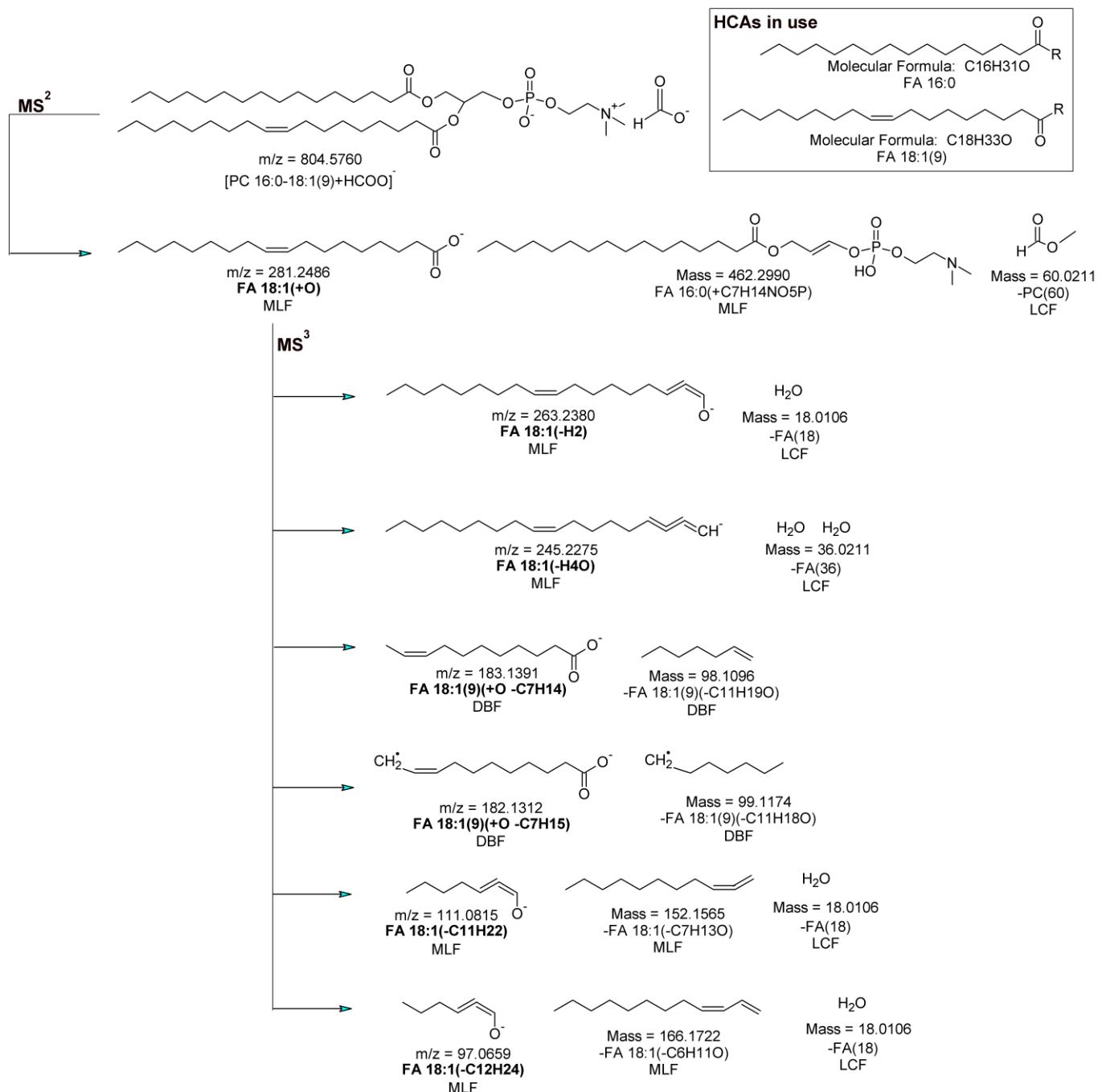


# S3D Fig) CL 14:1-14:1-14:1-15:1 (cont.)



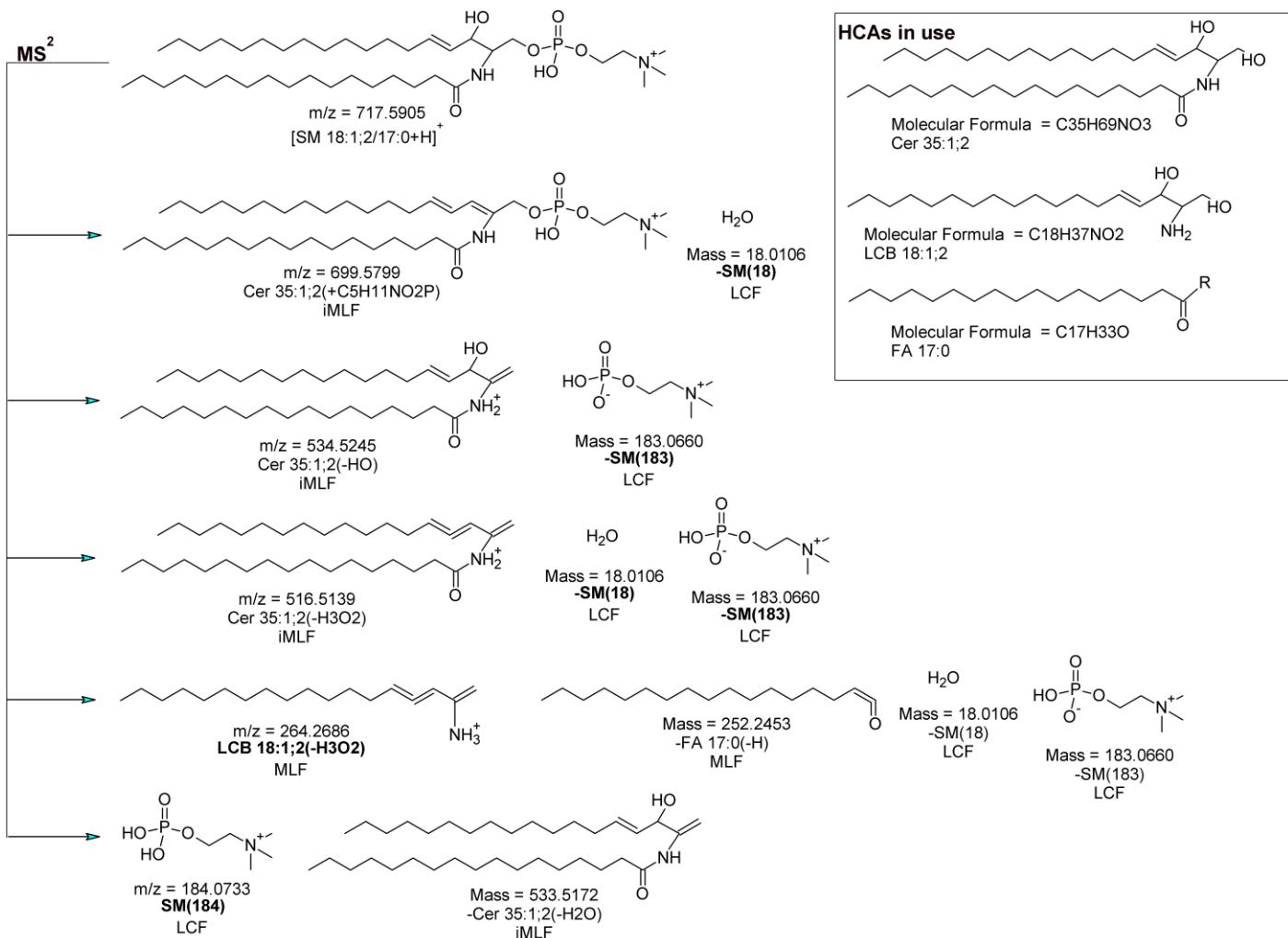
# S3E Fig) PC 16:0-18:1(9)

Proposed structures corresponding to spectrum shown in Fig. 3E.



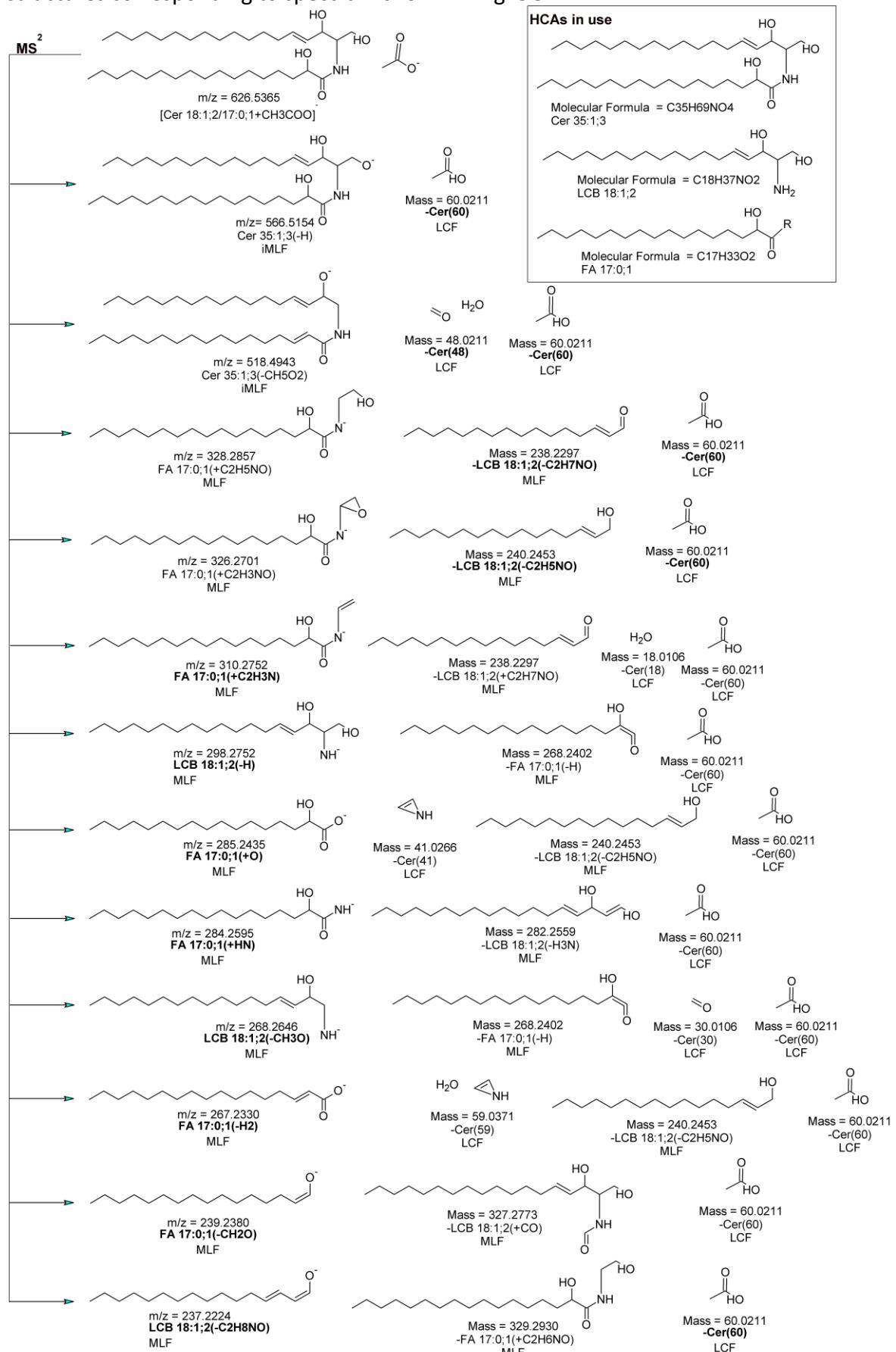
# S3F Fig) SM 18:1;2/17:0

Proposed structures corresponding to spectrum shown in Fig. 3F.



# S3G Fig) Cer 18:1;2/17:0;1

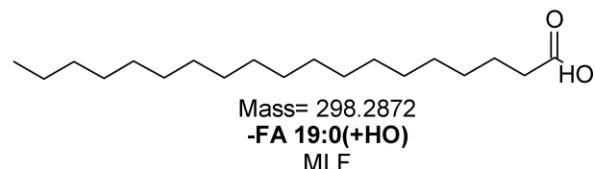
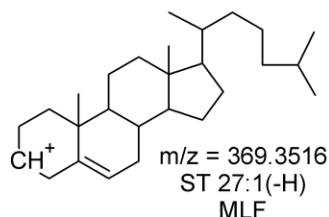
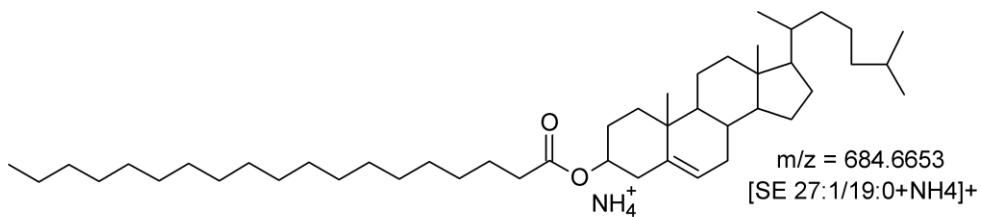
Proposed structures corresponding to spectrum shown in Fig. 3G.



# S3H Fig) SE 27:1/19:0

Proposed structures corresponding to spectrum shown in Fig. 3H.

MS<sup>2</sup>



NH<sub>3</sub>  
Mass = 17.0266  
-SE(17)  
LCF

