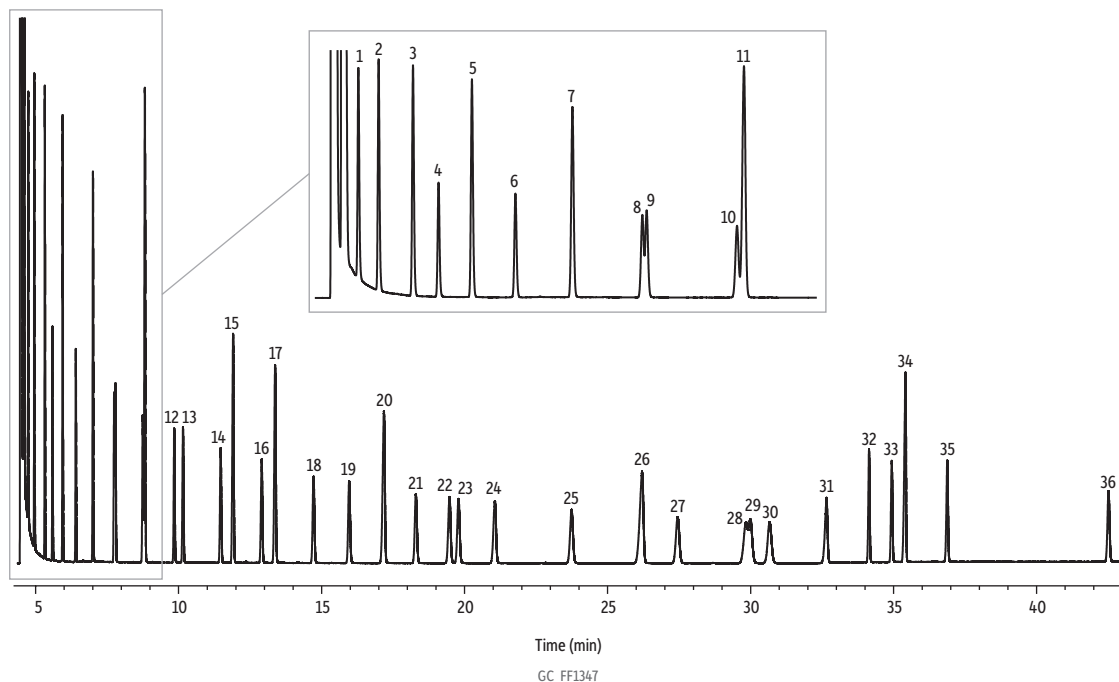


# Food Industry FAMES on Rt-2560 by AOCs Method Ce-1j-07 using H<sub>2</sub> (Original Method)



| Peaks                        | t <sub>r</sub> (min) | Conc. (µg/mL) | Structural Nomenclature       |
|------------------------------|----------------------|---------------|-------------------------------|
| 1. Methyl caproate           | 4.741                | 400           | C6:0                          |
| 2. Methyl octanoate          | 4.954                | 400           | C8:0                          |
| 3. Methyl decanoate          | 5.319                | 400           | C10:0                         |
| 4. Methyl undecanoate        | 5.589                | 200           | C11:0                         |
| 5. Methyl dodecanoate        | 5.942                | 400           | C12:0                         |
| 6. Methyl tridecanoate       | 6.403                | 200           | C13:0                         |
| 7. Methyl myristate          | 7.005                | 400           | C14:0                         |
| 8. Methyl myristoleate       | 7.744                | 200           | C14:1 (c9)                    |
| 9. Methyl pentadecanoate     | 7.790                | 200           | C15:0                         |
| 10. Methyl pentadecenoate    | 8.745                | 200           | C15:1 (C10)                   |
| 11. Methyl palmitate         | 8.818                | 600           | C16:0                         |
| 12. Methyl palmitoleate      | 9.846                | 200           | C16:1 (c9)                    |
| 13. Methyl heptadecanoate    | 10.153               | 200           | C17:0                         |
| 14. Methyl heptadecenoate    | 11.469               | 200           | C17:1 (c10)                   |
| 15. Methyl stearate          | 11.905               | 400           | C18:0                         |
| 16. Methyl octadecenoate     | 12.903               | 200           | C18:1 (t9)                    |
| 17. Methyl oleate            | 13.372               | 400           | C18:1 (c9)                    |
| 18. Methyl linolelaidate     | 14.718               | 200           | C18:2 (t9,t12)                |
| 19. Methyl linoleate         | 15.963               | 200           | C18:2 (c9,c12)                |
| 20. Methyl arachidate        | 17.176               | 400           | C20:0                         |
| 21. Methyl linolenate        | 18.295               | 200           | C18:3 (c6,c9,c12)             |
| 22. Methyl eicosenoate       | 19.470               | 200           | C20:1 (c11)                   |
| 23. Methyl linolenate        | 19.787               | 200           | C18:3 (c9,c12,c15)            |
| 24. Methyl heneicosanoate    | 21.060               | 200           | C21:0                         |
| 25. Methyl eicosadienoate    | 23.742               | 200           | C20:2 (c11,c14)               |
| 26. Methyl behenate          | 26.204               | 400           | C22:0                         |
| 27. Methyl eicosatrienoate   | 27.448               | 200           | C20:3 (c8,c11,c14)            |
| 28. Methyl erucate           | 29.836               | 200           | C22:1 (c13)                   |
| 29. Methyl eicosatrienoate   | 29.996               | 200           | C20:3 (c11,c14,c17)           |
| 30. Methyl arachidonate      | 30.667               | 200           | C20:4 (c5,c8,c11,c14)         |
| 31. Methyl tricosanoate      | 32.652               | 200           | C23:0                         |
| 32. Methyl docosadienoate    | 34.142               | 200           | C22:2 (c13,c16)               |
| 33. Methyl eicosapentaenoate | 34.934               | 200           | C20:5 (c5,c8,c11,c14,c17)     |
| 34. Methyl lignocerate       | 35.417               | 400           | C24:0                         |
| 35. Methyl nervonate         | 36.878               | 200           | C24:1 (c15)                   |
| 36. Methyl docosahexaenoate  | 42.513               | 200           | C22:6 (c4,c7,c10,c13,c16,c19) |

**Column** Rt-2560, 100 m, 0.25 mm ID, 0.20 µm (cat.# 13198)  
**Sample** Food industry FAME mix (cat.# 35077)  
**Diluent:** Hexane  
**Conc.:** 10,000 µg/mL total concentration  
**Injection**  
**Inj. Vol.:** 1 µL split (split ratio 20:1)  
**Liner:** Topaz 4.0 mm ID Precision inlet liner w/wool (cat.# 23305)  
**Inj. Temp.:** 250 °C  
**Oven**  
**Oven Temp.:** 180 °C (hold 32 min) to 215 °C at 20 °C/min (hold 30.25 min)  
**Carrier Gas** H<sub>2</sub>, constant flow  
**Flow Rate:** 2.5 mL/min  
**Detector** FID @ 250 °C  
**Constant Column +**  
**Constant Make-up:** 52 mL/min  
**Hydrogen flow:** 40 mL/min  
**Air flow:** 400 mL/min  
**Data Rate:** 50 Hz  
**Instrument** Agilent 7890A GC  
**Notes** Hydrogen flow optimized to achieve effective linear velocity (www.restek.com/FAMEH2).  
 C4:0 Methyl butyrate (623-42-7) elutes in the solvent front.  
 Resolution of critical pair *cis*-11-C20:1 (#22) and *cis*-9,12,15-C18:3 (#23) is 2.03.