

## CORRIGENDUM

# 5-Hydroxy-5,6-dihydro-apo-4,4'-lycopene and methyl 5-hydroxy-5,6-dihydro-apo-4,4'-lycopenoate, novel C<sub>30</sub>-carotenoids produced by a mutant of marine bacterium *Halobacillus halophilus*

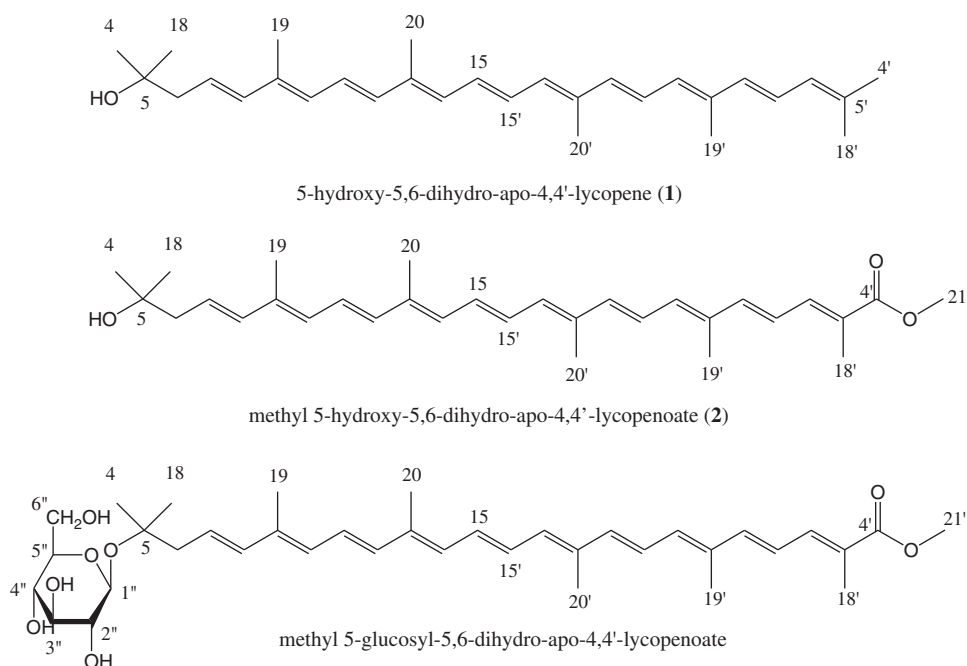
Ayako Osawa, Yoko Ishii, Nao Sasamura, Marie Morita, Saskia Köcher, Volker Müller, Gerhard Sandmann and Kazutoshi Shindo

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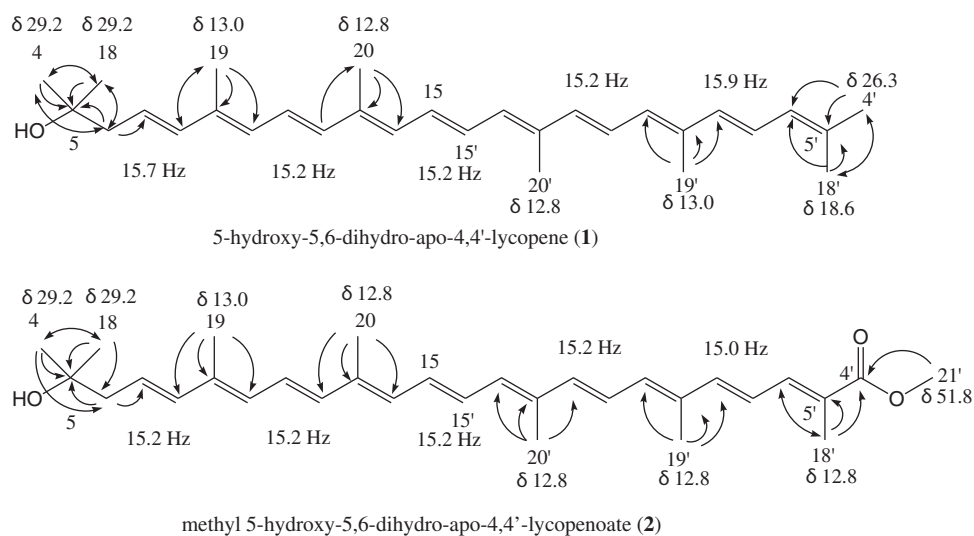
**Correction to:** *The Journal of Antibiotics* (2010) 63: 291–295; doi:10.1038/ja.2010.33; published online 23 April 2010

shown below. Reflecting this correction, title of the article has been changed. Authors would like to apologize for the inconvenience caused.

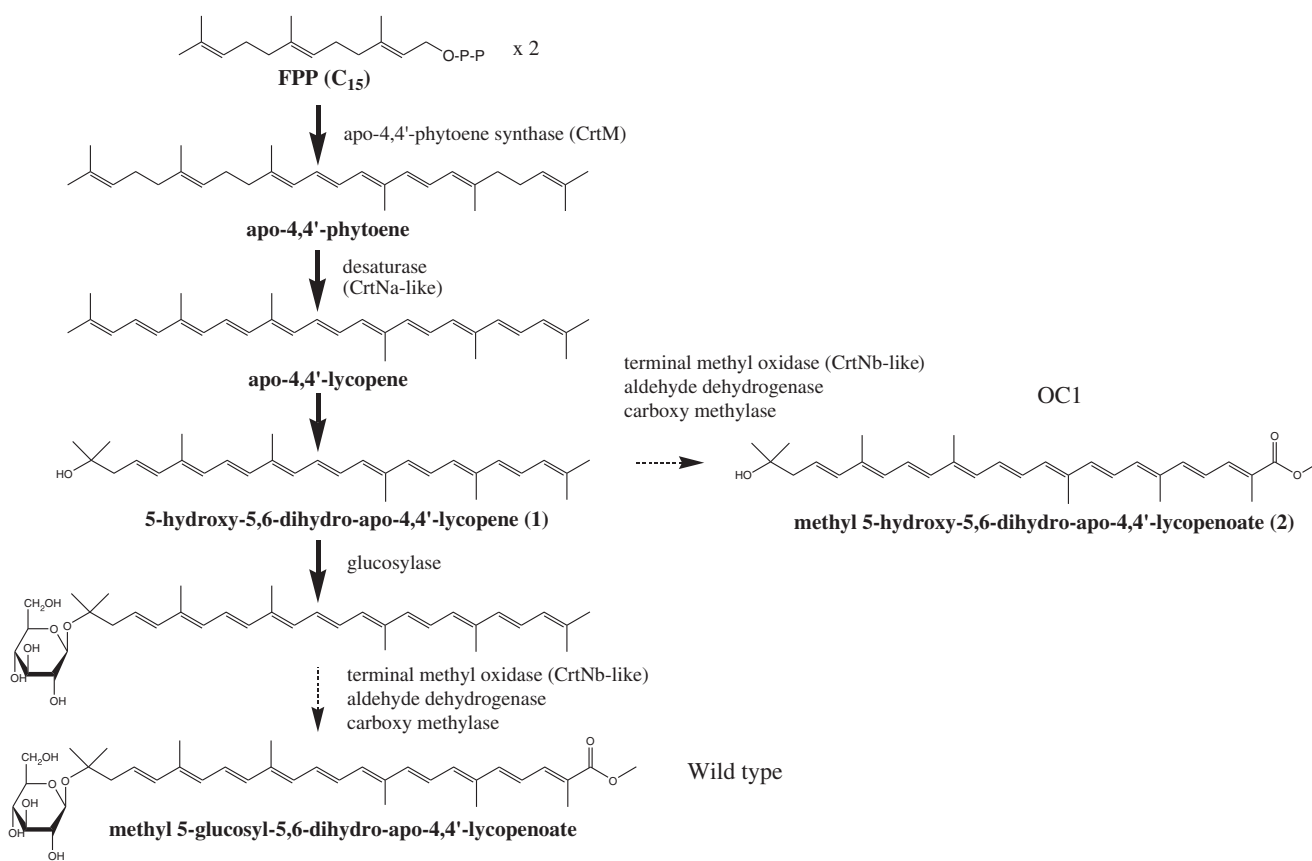
The authors of the above Article noted an error in publication of this paper in Figures 1–3 and Table 1. The corrected figures and table are



**Figure 1** The structure of carotenoids **1** and **2** produced by mutant OC1 (a mutant of *H. halophilus*), and methyl 5-glucosyl-5,6-dihydro-apo-4,4'-lycopenoate produced by *H. halophilus*.



**Figure 2** Key  $^1\text{H}$ - $^{13}\text{C}$  long-range couplings,  $J$  values, and  $\delta_c$  values observed in the NMR analyses of **1** and **2**.



**Figure 3** Proposed biosynthetic pathway of methyl 5-glucosyl-5,6-dihydro-apo-4,4'-lycopenoate in *H. halophilus* DSM2266.

**Table 1**  $^1\text{H}$  and  $^{13}\text{C}$  NMR data for 5-hydroxy-5,6-dihydro-apo-4,4'-lycopene (**1**), methyl 5-hydroxy-5,6-dihydro-apo-4,4'-lycopenoate (**2**) and **2** peracetate [**9**] in  $\text{CDCl}_3$ 

Position	<b>1</b>		<b>2</b>		<b>2 Peracetate</b>	
	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$	$\delta_{\text{H}}$	$\delta_{\text{C}}$
4	1.24 (3H, s)	29.2 ( $\text{CH}_3$ )	1.24 (3H, s)	29.3 ( $\text{CH}_3$ )	1.17 (3H, s)	26.6 ( $\text{CH}_3$ )
5		71.0 (C)		71.0 (C)		78.8 (C)
6	2.31 (2H, d, 7.4)	47.5 ( $\text{CH}_2$ )	2.32 (2H, d, 7.4)	47.5 ( $\text{CH}_2$ )	2.24 (1H, dd, 7.9, 13.8) 2.42 (1H, dd, 7.4, 13.8)	45.9 ( $\text{CH}_2$ )
7	5.76 (1H, ddd, 7.4, 7.4, 15.7)	124.5 (CH)	5.78 (1H, ddd, 7.4, 7.4, 15.7)	124.6 (CH)	5.72 (1H, ddd, 7.4, 7.9, 15.1)	125.2 (CH)
8	6.21 (1H, d, 15.7)	138.8 (CH)	6.21 (1H, d, 15.2)	138.8 (CH)	6.13 (1H, d, 15.1)	137.9 (CH)
9		135.0 (C)		135.0 (C)		135.5 (C)
10	6.14 (1H, d, 11.4)	131.2 (CH)	6.14 (1H, d, 10.9)	130.9 (CH)	6.08 (1H, d, 11.3)	130.9 (CH)
11	6.61 (1H, dd, 11.4, 15.2)	124.9 (CH)	6.62 (1H, dd, 10.9, 15.2)	125.1 (CH)	6.63 (1H, dd, 11.3, 14.4)	125.0 (CH)
12	6.35 (1H, d, 15.2)	137.8 (CH)	6.36 (1H, d, 15.2)	137.6 (CH)	6.35 (1H, d, 14.4)	137.4 (CH)
13		136.2 (C) <sup>a</sup>		136.3 (C) <sup>a</sup>		136.0 (C) <sup>a</sup>
14	6.25 (1H, d, 10.3)	132.7 (CH) <sup>a</sup>	6.24 (1H, d, 10.3)	132.5 (CH)	6.26 (1H, d, 9.7)	132.5 (CH)
15	6.63 (1H, dd, 10.3, 15.2)	130.0 (CH) <sup>a</sup>	6.65 (1H, dd, 10.3, 15.2)	131.2 (CH)	6.65 (1H, dd, 9.7, 15.2)	130.8 (CH)
18	1.24 (3H, s)	29.2 ( $\text{CH}_3$ )	1.26 (3H, s)	29.5 ( $\text{CH}_3$ )	1.19 (3H, s)	25.6 ( $\text{CH}_3$ )
19	1.93 (3H, s)	13.0 ( $\text{CH}_3$ )	1.94 (3H, s)	13.0 ( $\text{CH}_3$ )	1.91 (3H, s)	13.0 ( $\text{CH}_3$ )
20	1.97 (3H, s)	12.8 ( $\text{CH}_3$ )	1.98 (3H, s)	12.8 ( $\text{CH}_3$ )	1.98 (3H, s)	12.8 ( $\text{CH}_3$ )
4'	1.82 (3H, s)	26.3 ( $\text{CH}_3$ )		169.0 (C)		169.1 (C)
5'		135.8 (C)		125.7 (C)		125.7 (C)
6'	5.94 (1H, d, 10.8)	126.1 (CH)	7.30 (1H, d, 12.3)	139.1 (CH)	7.30 (1H, d, 11.3)	139.1 (CH)
7'	6.47 (1H, dd, 10.8, 15.9)	124.7 (CH)	6.52 (1H, dd, 12.3, 15.0)	122.9 (CH)	6.51 (1H, dd, 11.3, 15.1)	122.9 (CH)
8'	6.21 (1H, d, 15.9)	135.0 (CH)	6.66 (1H, d, 15.0)	144.2 (CH)	6.62 (1H, d, 15.1)	144.2 (CH)
9'		136.6 (C) <sup>a</sup>		135.2 (C) <sup>a</sup>		135.1 (C)
10'	6.18 (1H, d, 12.0)	131.4 (CH)	6.32 (1H, d, 11.9)	136.2 (CH)	6.35 (1H, d, 11.4)	136.2 (CH)
11'	6.63 (1H, m)	125.2 (CH)	6.63 (1H, dd, 11.9, 15.2)	125.1 (CH)	6.63 (1H, dd, 11.4, 15.0)	125.0 (CH)
12'	6.35 (1H, d, 15.2)	137.3 (CH)	6.45 (1H, d, 15.2)	139.9 (CH)	6.45 (1H, d, 15.0)	139.9 (CH)
13'		136.3 (C) <sup>a</sup>		136.0 (C)		137.0 (C) <sup>a</sup>
14'	6.25 (1H, d, 12.0)	132.5 (CH) <sup>a</sup>	6.29 (1H, d, 11.0)	134.1 (CH)	6.31 (1H, d, 11.4)	134.1 (CH)
15'	6.63 (1H, dd, 12.0, 15.2)	130.2 (CH) <sup>a</sup>	6.63 (1H, dd, 11.0, 15.2)	129.9 (CH)	6.63 (1H, dd, 11.4, 15.2)	129.9 (CH)
18'	1.82 (3H, s)	18.6 ( $\text{CH}_3$ )	2.00 (3H, s)	12.8 ( $\text{CH}_3$ )	2.00 (3H, s)	12.8 ( $\text{CH}_3$ )
19'	1.97 (3H, s)	13.0 ( $\text{CH}_3$ )	2.00 (3H, s)	12.8 ( $\text{CH}_3$ )	2.00 (3H, s)	12.8 ( $\text{CH}_3$ )
20'	1.97 (3H, s)	12.8 ( $\text{CH}_3$ )	2.00 (3H, s)	12.8 ( $\text{CH}_3$ )	1.98 (3H, s)	12.8 ( $\text{CH}_3$ )
$\text{COOCH}_3$			3.77 (3H, s)	51.8 ( $\text{CH}_3$ )	3.77 (3H, s)	51.7 ( $\text{CH}_3$ )
1''					4.69 (1H, d, 7.8)	93.4 (CH)
2''					4.98 (1H, dd, 7.8, 9.5)	71.6 (CH)
3''					5.24 (1H, dd, 9.0, 9.5)	73.1 (CH)
4''					5.04 (1H, dd, 9.0, 10.0)	68.9 (CH)
5''					3.68 (1H, ddd, 1.5, 5.8, 10.0)	71.6 (CH)
6''					4.10 (1H, dd, 1.5, 12.1) 4.22 (1H, dd, 5.8, 12.1) 4.22 (1H, dd, 5.8, 12.1)	62.5 ( $\text{CH}_2$ )

<sup>13</sup>C signals of acetyl groups in **2** peracetate were observed at  $\delta = 20.6\text{--}20.7$  ( $\text{CH}_3$ ) and  $\delta = 169.0\text{--}170.6$  (C=O).

<sup>a</sup>Interchangeable.