Table S2. NMR Assignments for *In Vivo* Produced Methylated Triterpenes [1]

C31a botryococcene by TMT3	
C positions	δ
1	17.78
2	131.37
3	124.47
4	26.82
5	39.81
6	134.76
7	124.91
8	23.20
9	41.41
10	42.12
11	135.91
12	133.82
13	36.84
14	37.51
15	25.92
16	124.58
17	135.11
18	37.62
19	33.44
20	40.82
21	150.27
22	109.44
23	25.79
24	16.00
25	23.63
26	146.85
27	111.20
28	21.23
29	16.07
30	19.06
31	19.76

C31b botryococcene by TMT3	
C positions	δ
1	109.44
2	150.24
3	40.82
4	33.44
5	37.58
6	135.07
7	124.80
8	23.20
9	41.45
10	42.12
11	135.89
12	133.84
13	36.77
14	37.46
15	25.90
16	124.69
17	134.80
18	39.81
19	26.82
20	124.50
21	131.34
22	17.78
23	19.06
24	16.00
25	23.63
26	146.85
27	111.20
28	21.23
29	16.07
30	25.79
31	19.76

C32 botryococcene by TMT3	
C positions	δ
1	109.44
2	150.23
3	40.82
4	33.44
5	37.58
6	135.07
7	124.70
8	23.19
9	41.45
10	42.11
11	135.91
12	133.82
13	36.84
14	37.51
15	25.92
16	124.58
17	135.11
18	37.62
19	33.42
20	40.84
21	150.23
22	109.44
23	19.04
24	16.00
25	23.62
26	146.84
27	111.19
28	21.23
29	16.07
30	19.04
31	19.78
32	19.78

C31 squalene by TMT1	
C positions	δ
1	109.46
2	150.25
3	40.79
4	33.42
5	37.59
6	134.99a
7	124.49b
8	26.85
9	39.84
10	135.19a
11	124.40b
12	28.36
13	28.36
14	124.40b
15	135.17a
16	39.81c
17	26.74d
18	124.36
19	135.25a
20	39.81c
21	26.72d
22	124.19
23	131.34
24	17.76
25	19.03
26	16.09e
27	16.12e
28	16.12e
29	16.09e
30	25.79
31	19.76

C32 squalene by TMT1	
C positions	δ
1	109.46
2	150.23
3	40.79
4	33.43
5	37.61
6	135.15f
7	124.19
8	26.73
9	39.86
10	135.25f
11	124.41
12	28.37
13	28.37
14	124.41
15	135.25f
16	39.86
17	26.73
18	124.19
19	135.15f
20	37.61
21	33.43
22	40.79
23	150.23
24	109.46
25	19.03
26	16.08
27	16.12
28	16.12
29	16.08
30	19.03
31	19.78
32	19.78

C31 squalene by TMT2	
C positions	δ
1	109.46
2	150.26
3	40.79
4	33.43
5	37.59
6	134.99a
7	124.49b
8	26.85
9	39.84
10	135.19a
11	124.40b
12	28.36
13	28.36
14	124.40b
15	135.17a
16	39.81c
17	26.74d
18	124.36
19	135.25a
20	39.81c
21	26.72d
22	124.19
23	131.34
24	17.76
25	19.03
26	16.08e
27	16.13e
28	16.13e
29	16.08e
30	25.79
31	19.76

C32 squalene by TMT2	
C positions	δ
1	109.45
2	150.26
3	40.79
4	33.42
5	37.59
6	135.17f
7	124.18
8	26.71
9	39.86
10	135.25f
11	124.40
12	28.35
13	28.35
14	124.40
15	135.25f
16	39.86
17	26.71
18	124.18
19	135.17f
20	37.59
21	33.42
22	40.79
23	150.26
24	109.45
25	19.04
26	16.08g
27	16.12g
28	16.12g
29	16.08g
30	19.04
31	19.76
32	19.76

C31 squalene by TMT3	
C positions	δ
1	109.45
2	150.25
3	40.79
4	33.43
5	37.60
6	134.99a
7	124.49b
8	26.86
9	39.84
10	135.19a
11	124.40b
12	28.36
13	28.36
14	124.40b
15	135.17a
16	39.81c
17	26.75d
18	124.36
19	135.25a
20	39.81c
21	26.72d
22	124.19
23	131.34
24	17.75
25	19.03
26	16.09e
27	16.12e
28	16.12e
29	16.09e
30	25.79
31	19.76

C32 squalene by TMT3	
C positions	δ
1	109.46
2	150.25
3	40.79
4	33.42
5	37.60
6	135.16f
7	124.18
8	26.73
9	39.86
10	135.26f
11	124.40
12	28.36
13	28.36
14	124.40
15	135.26f
16	39.86
17	26.73
18	124.18
19	135.16f
20	37.60
21	33.42
22	40.79
23	150.25
24	109.46
25	19.04
26	16.08g
27	16.12g
28	16.12g
29	16.08g
30	19.04
31	19.78
32	19.78

a, b, c, d, e, f, g: Signals with the same letter may be interchangeable in a same column.

^{1}H and ^{13}C NMR spectra were recorded on a JEOL alpha 600 NMR spectrometer at 300K. Chemical shifts were referenced relative to solvent peaks, namely d_H 7.24 and d_C 77.1 for $CDCl_3$. Each product was identified as shown in Fig. 4 by referring ^{13}C chemical shifts for botryococcenes and methylsqualenes to those previously reported (1-4).

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