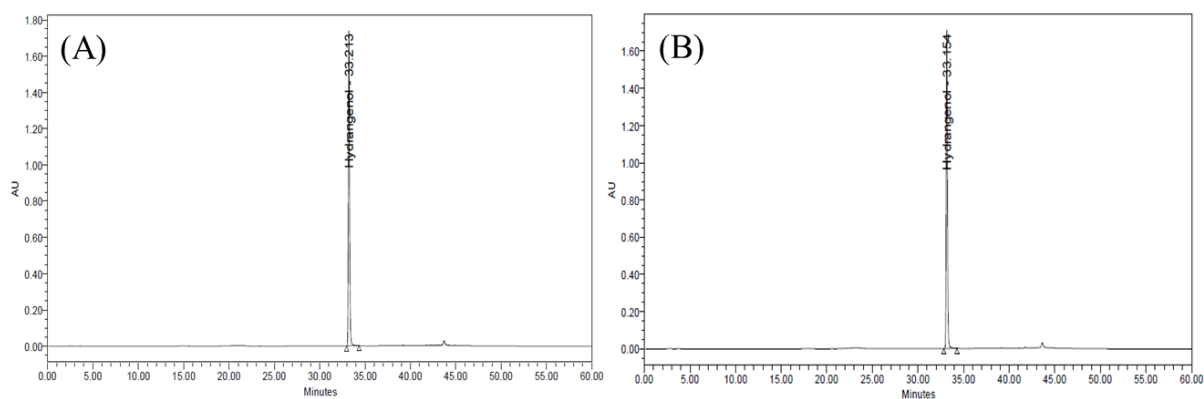


**Table S1.** The primer sequences for qRT-PCR

<b>Gene</b>	<b>Forward sequence</b>	<b>Reverse sequence</b>
<b>HYAL1</b>	5' TACACAGCATGCTCAGAAAG 3'	5' AGTGTCTCCATTCCAAACAG 3'
<b>HYAL2</b>	5' AATCTGTGGAACGCTACATC 3'	5' GCCTGCTTCATTACTCTGTC 3'
<b>Involucrin</b>	5' GGGTCAGTCACTTAAGCAAG 3'	5' CTACTTCTCCTGCTGTGTCC 3'
<b>Filaggrin</b>	5' CTCAGGAGGAAGAGGACAGT 3'	5' CAAGGTGCTTTGCTGTAAAT 3'
<b>COL1A1</b>	5' CCCAGAACATCACCTATCAC 3'	5' GAGGTCTTGGTGGTTTTGTA 3'
<b>MMP-1</b>	5' TTGCCCAGAGAAAAGCTTCAG 3'	5' TAGCAGCCCAGAGAAGCAACA 3'
<b>MMP-3</b>	5' GTTCTGGGCTATACGAGGGC 3'	5' GGCAGCATCGATCTTCTTCA 3'
<b>COX-2</b>	5' TGCTGTACAAGCAGTGGCAA 3'	5' GCAGCCATTTCCTTCTCTCC 3'
<b>IL-6</b>	5' GAGGATACTCACTCCCAACAGACC 3'	5' AAGTGCATCATCGTTGTTCATACA 3'
<b>NQO-1</b>	5' TGGCCGATTCAGAGTGGCAT 3'	5' AAACAGGCTGCTTGGAGCAAAA 3'
<b>GCLM</b>	5' GGGAACCTGCTCAACTGGGG 3'	5' CTGCATGGGCATGGTGCATT 3'
<b>GCLC</b>	5' TCCGGCATCGGAGAGGAGA 3'	5' AGCAGTTGCCCATCCCGAAT 3'
<b><math>\beta</math>-actin</b>	5' ATCACTATTGGCAACGAGCG 3'	5' TCAGCAATGCCTGGGTACAT 3'

**Figure S1.**



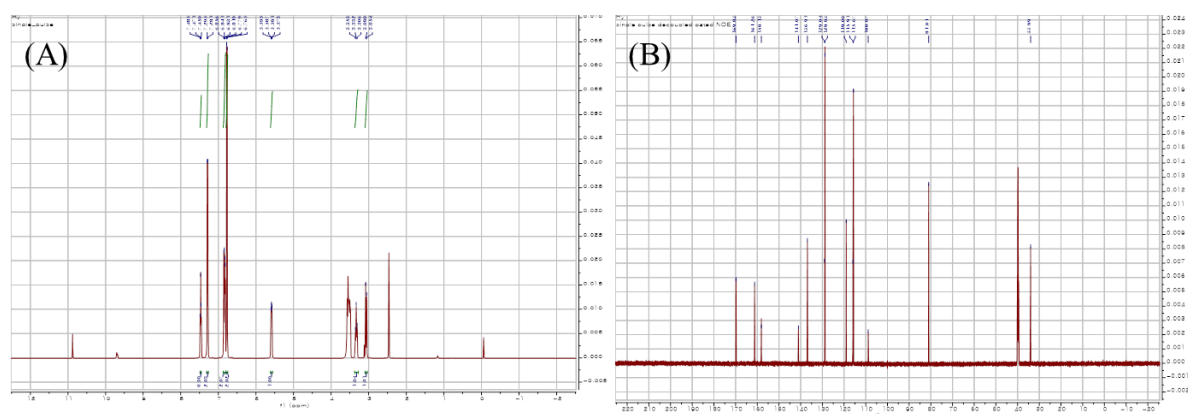
<HPLC analysis conditions>

- Column : Luna C18 (5  $\mu$ m, 250  $\times$  4.6 mm, Phenomenex)
- Detector : Waters 2489 UV/Vis Detector (228 nm)
- Pump : Waters e2695
- Oven temperature : 30°C
- Injection volume : 20  $\mu$ L
- Flow rate : 1 mL/min
- Solvent system : Gradient system (linear gradient system)

Time (min)	% A (ACN)	% B (H <sub>2</sub> O)
<b>Initial</b>	20	80
<b>15</b>	25	75
<b>30</b>	50	50
<b>40</b>	100	0
<b>50</b>	20	80

**Figure S1.** (A) HPLC profile of hydrangenol isolated from WHS and (B) authentic reference hydrangenol

**Figure S2.**



<sup>1</sup>H-NMR (600 MHz, DMSO-*d*<sub>6</sub>-D<sub>2</sub>O)  
δ: 7.47 (1H, dd, *J* = 7.2, 8.4 Hz, H-6), 7.29 (2H, d, *J* = 8.4 Hz, H-2',6'), 6.85 (1H, d, *J* = 8.4 Hz, H-7), 6.83 (1H, d, *J* = 7.2 Hz, H-5), 6.77 (2H, d, *J* = 8.4 Hz, H-3',5'), 5.50 (1H, dd, *J* = 3.6, 12.0 Hz, H-3), 3.33 (1H, dd, *J* = 12.0, 15.6 Hz, H-4a), 3.07 (1H, dd, *J* = 3.6, 15.6 Hz, H-4b).

<sup>13</sup>C-NMR (125 MHz, DMSO-*d*<sub>6</sub>-D<sub>2</sub>O)  
δ: 34.0 (C-4), 81.0 (C-3), 108.9 (C-9), 115.7 (C-3',5'), 115.9 (C-7), 119.0 (C-5), 128.8 (C-2',6'), 129.0 (C-1'), 136.9 (C-6), 141.1 (C-10), 158.1 (C-4'), 161.3 (C-8), 169.8 (C-1).

**Figure S2.** (A) <sup>1</sup>H and (B) <sup>13</sup>C NMR spectra of hydrangenol isolated from WHS measured in DMSO-*d*<sub>6</sub>-H<sub>2</sub>O. The chemical structure of hydrangenol was elucidated by instrumental analyses. 1D-NMR experiments were conducted using (A) <sup>1</sup>H-(600 MHz) and (B) <sup>13</sup>C-NMR (150 MHz), and recorded with VNS (JEOL, Tokyo, Japan).