



**Molecular Formula:** C<sub>18</sub>H<sub>18</sub>O<sub>7</sub>  
**Molecular Weight:** 346.336  
**Accurate Mass:** 346.105255  
**Percentage Composition:** C 62.42%; H 5.24%; O 32.34%  
**Biological Source:** Constit. of the roots of *Peucedanum ostruthium*  
**Physical Description:** Needles (MeOH)  
**Melting Point:** Mp 116-119°

>> **Derivative:** 2'-*O*-(2-Hydroxypropanoyl)  
**Synonym(s):** Fesumtuorin A

**Molecular Formula:** C<sub>19</sub>H<sub>20</sub>O<sub>8</sub>  
**Molecular Weight:** 376.362  
**Accurate Mass:** 376.11582  
**Percentage Composition:** C 60.64%; H 5.36%; O 34.01%  
**Biological Source:** Constit. of the dried roots of *Ferula sumbul*  
**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>25</sup> +17.4 (c, 1.1 in MeOH)  
**UV:** [neutral]  $\lambda_{\max}$  220 (log  $\epsilon$  4.24); 250 (log  $\epsilon$  4.07); 267 (log  $\epsilon$  4.01); 311 (log  $\epsilon$  3.92) (MeOH)  
**Other Data:** Stereochem. not determined

>> **Derivative:** 2'-*O*-(3-Methylbutanoyl)

**CAS Registry Number:** 68006-84-8

**Molecular Formula:** C<sub>21</sub>H<sub>24</sub>O<sub>7</sub>  
**Molecular Weight:** 388.416  
**Accurate Mass:** 388.152205  
**Percentage Composition:** C 64.94%; H 6.23%; O 28.83%  
**Biological Source:** Constit. of *Ammi majus*  
**Physical Description:** Cryst.  
**Melting Point:** Mp 109-110°  
**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>25</sup> -26.4 (CHCl<sub>3</sub>)

>> **Derivative:** 2'-*O*-(3-Methyl-2-butenoyl)  
**Synonym(s):** Senecioplprangol

**Molecular Formula:** C<sub>21</sub>H<sub>22</sub>O<sub>7</sub>  
**Molecular Weight:** 386.401  
**Accurate Mass:** 386.136555  
**Percentage Composition:** C 65.28%; H 5.74%; O 28.98%  
**Biological Source:** Constit. of *Ferulago capillaris*  
**Physical Description:** Cryst. (MeOH)  
**Melting Point:** Mp 164-165°  
**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>20</sup> +9 (c, 1 in Me<sub>2</sub>CO)  
**UV:** [neutral]  $\lambda_{\max}$  220 (log  $\epsilon$  4.1); 250 (log  $\epsilon$  3.83); 258 (log  $\epsilon$  3.7); 267 (log  $\epsilon$  3.7); 309 (log  $\epsilon$  3.6) (EtOH)

>> **Derivative:** 2'-Angeloyl  
**Synonym(s):** Ostruthol

**CAS Registry Number:** 642-08-0

**Molecular Formula:** C<sub>21</sub>H<sub>22</sub>O<sub>7</sub>  
**Molecular Weight:** 386.401  
**Accurate Mass:** 386.136555  
**Percentage Composition:** C 65.28%; H 5.74%; O 28.98%  
**Biological Source:** Isol. from *Ammi pachycarpa*, *Ammi archangelica*, *Peucedanum ostruthium* and *Peucedanum hispanicum*  
**Physical Description:** Cryst. (C<sub>6</sub>H<sub>6</sub>)  
**Melting Point:** Mp 136-137°  
**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>15</sup> -18.3 (Py)