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|--------------------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Product Name</b>      | : Fenretinide                                                                                                                                                                      |
| <b>Synonyms</b>          | : 4-HPR; 4-Hydroxy(phenyl)retinamide; MK-4016; Retinoic Acid p-hydroxyphenylamide; Ro 22-4667                                                                                      |
| <b>Cat No.</b>           | : M15496                                                                                                                                                                           |
| <b>CAS Number</b>        | : 65646-68-6                                                                                                                                                                       |
| <b>Molecular Formula</b> | : C <sub>26</sub> H <sub>33</sub> NO <sub>2</sub>                                                                                                                                  |
| <b>Formula Weight</b>    | : 391.55                                                                                                                                                                           |
| <b>Chemical Name</b>     | : (2E,4E,6E,8E)-N-(4-hydroxyphenyl)-3,7-dimethyl-9-(2,6,6-trimethylcyclohex-1-en-1-yl)nona-2,4,6,8-tetraenamide                                                                    |
| <b>Description</b>       | : Fenretinide (4-HPR) is a synthetic retinoid derivative. 4-HBR is shown to exhibit binding to the retinoic acid receptors (RAR) at concentrations necessary to induce cell death. |
| <b>Pathway</b>           | : Metabolic Enzyme/Protease                                                                                                                                                        |
| <b>Target</b>            | : Retinoid Receptor                                                                                                                                                                |
| <b>Receptor</b>          | : RAR/RAX                                                                                                                                                                          |
| <b>Solubility</b>        | : DMSO: ≥100mM                                                                                                                                                                     |
| <b>SMILES</b>            | : <chem>O=C(NC1=CC=C(O)C=C1)/C=C(C)/C=C/C=C(C)/C=C/C2=C(C)CCCC2(C)C</chem>                                                                                                         |
| <b>Storage</b>           | : (-20°C)                                                                                                                                                                          |
| <b>Stability</b>         | : ≥ 2 years                                                                                                                                                                        |
| <b>Reference</b>         | :                                                                                                                                                                                  |

1. Manzano VM, et al. J Pharmacol Exp Ther, 1999, 289(1), 123-132.