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| <b>Product Name</b>      | : Fexaramine                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| <b>Synonyms</b>          | : Fexaramine                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| <b>Cat No.</b>           | : M15095                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| <b>CAS Number</b>        | : 574013-66-4                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| <b>Molecular Formula</b> | : C32H36N2O3                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| <b>Formula Weight</b>    | : 496.64                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| <b>Chemical Name</b>     | : 2-Propenoic acid, 3-[3-[(cyclohexylcarbonyl)[[4'-(dimethylamino)[1,1'-biphenyl]-4-yl]methyl]amino]phenyl]-, methyl ester                                                                                                                                                                                                                                                                                                                            |
| <b>Description</b>       | : A potent, specific FXR agonist with EC50 of 25 nM, more efficacious ligand for FXR than GW4064; shows no transcriptional activity against hRXR $\alpha$ , hPPAR $\alpha$ / $\gamma$ / $\delta$ , mPXR, hPXR, hLXR $\alpha$ , hTR $\beta$ , hRAR $\beta$ , mCAR, mERR $\gamma$ , and hVDR; strongly induces target genes SHP, PLTP BSEP, and MRP-2 in the HEPG2 liver cell system; robustly induces FGF15, leading to alterations in BA composition. |
| <b>Pathway</b>           | : Metabolic Enzyme/Protease                                                                                                                                                                                                                                                                                                                                                                                                                           |
| <b>Target</b>            | : FXR                                                                                                                                                                                                                                                                                                                                                                                                                                                 |
| <b>Receptor</b>          | : FXR                                                                                                                                                                                                                                                                                                                                                                                                                                                 |
| <b>Solubility</b>        | : 10 mM in DMSO                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| <b>SMILES</b>            | : <chem>OC(=O)C=C/C1=CC=CC(N(C(C2CCCCC2)=O)CC3=CC=C(C4=CC=C(N(C)C)C=C4)C=C3)=C1</chem>                                                                                                                                                                                                                                                                                                                                                                |
| <b>Storage</b>           | : (-20°C)                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| <b>Stability</b>         | : $\geq 2$ years                                                                                                                                                                                                                                                                                                                                                                                                                                      |
| <b>Reference</b>         | :                                                                                                                                                                                                                                                                                                                                                                                                                                                     |

1. Downes M, et al. Mol Cell. 2003 Apr;11(4):1079-92. | 2. Fang S, et al. Nat Med. 2015 Feb;21(2):159-65. | 3. Cho SW, et al. J Bone Miner Res. 2013 Oct;28(10):2109-21.