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| <b>Product Name</b>      | : Acrizanib                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| <b>Synonyms</b>          | : LHA510;LHA 510                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| <b>Cat No.</b>           | : M10899                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
| <b>CAS Number</b>        | : 1229453-99-9                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |
| <b>Molecular Formula</b> | : C <sub>20</sub> H <sub>18</sub> F <sub>3</sub> N <sub>7</sub> O <sub>2</sub>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |
| <b>Formula Weight</b>    | : 445.41                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
| <b>Chemical Name</b>     | : 5-{[6-[(methylamino)methyl]pyrimidin-4-yl]oxy}-N-[1-methyl-5-(trifluoromethyl)-1H-pyrazol-3-yl]-1H-indole-1-carboxamide                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| <b>Description</b>       | <p>Acrizanib (LHA510, LHA-510) is a novel potent, selective VEGFR-2 (KDR) inhibitor with IC<sub>50</sub> of 17.4 nM in cell-based assays; exhibits &lt;10% remaining kinase activity against only 13 wildtype kinases, CSF1R, Kit, PDGFR<math>\alpha</math>, PDGFR<math>\beta</math>, VEGFR1, VEGFR2, VEGFR3, Fms, DDR1, DDR2, TIE1, and ABL1 in a panel of &gt;400 kinases; demonstrates potency and efficacy in rodent CNV models, limited systemic exposure after topical ocular administration, multiple formulation options, and an acceptable rabbit ocular PK profile. Other Indication Phase 1 Clinical</p> |
| <b>Pathway</b>           | : Angiogenesis                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      |
| <b>Target</b>            | : VEGFR                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| <b>Receptor</b>          | : VEGFR                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             |
| <b>Solubility</b>        | : 10 mM in DMSO                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |
| <b>SMILES</b>            | : <chem>O=C(N1C=CC2=C1C=CC(OC3=NC=NC(CNC)=C3)=C2)NC4=NN(C)C(C(F)(F)F)=C4</chem>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                     |
| <b>Storage</b>           | : (-20°C)                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| <b>Stability</b>         | : $\geq 2$ years                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| <b>Reference</b>         | :                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                   |

1.Adams CM, et al. J Med Chem. 2018 Feb 5. doi: 10.1021/acs.jmedchem.7b01731.