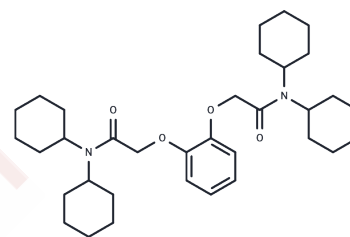


## Sodium ionophore III

## Chemical Properties

CAS No. : 81686-22-8  
 Formula: C<sub>34</sub>H<sub>52</sub>N<sub>2</sub>O<sub>4</sub>  
 Molecular Weight: 552.79  
 Appearance: no data available  
 Storage: Powder: -20°C for 3 years | In solvent: -80°C for 1 year



## Biological Description

|               |   |
|---------------|---|
| Description   | Sodium ionophore III (ETH2120) is a Na <sup>+</sup> ionophore ideal for testing sodium activity in plasma, serum, and blood.  |
| Targets(IC50) | Sodium Channel  |
| In vitro      | Preincubation of the cells with Sodium ionophore III not only completely abolished ATP synthesis but stimulated caffeate reduction. The addition of Sodium ionophore III to cells in the steady state of caffeate reduction dissipated the intracellular ATP level[1]. Lactate-sulfate grown cells are insensitive to Sodium ionophore III. The Sodium ionophore III ligand is a very effective receptor for the Am <sup>3+</sup> and Eu <sup>3+</sup> cations. |

## Solubility Information

|            |   |
|------------|---|
| Solubility | DMSO: Slightly soluble,<br>( $< 1$ mg/ml refers to the product slightly soluble or insoluble) |
|------------|---|

## Preparing Stock Solutions

|       | 1mg       | 5mg       | 10mg       |
|-------|-----------|-----------|------------|
| 1 mM  | 1.809 mL  | 9.045 mL  | 18.0901 mL |
| 5 mM  | 0.3618 mL | 1.809 mL  | 3.618 mL   |
| 10 mM | 0.1809 mL | 0.9045 mL | 1.809 mL   |
| 50 mM | 0.0362 mL | 0.1809 mL | 0.3618 mL  |

Please select the appropriate solvent to prepare the stock solution, according to the solubility of the product in different solvents. Please use it as soon as possible.

## Reference

- Imkamp F, et al. Chemiosmotic energy conservation with Na<sup>+</sup> as the coupling ion during hydrogen-dependent caffeate reduction by *Acetobacterium woodii*. J Bacteriol. 2002 Apr;184(7):1947-51.
- Wang L, et al. The role of Rnf in ion gradient formation in *Desulfovibrio alaskensis*. PeerJ. 2016 Apr 14;4:e1919.
- Makrlík, E., Kvíčalová, M. & Vaňura, P. J Solution Chem (2016) 45: 463. <https://doi.org/10.1007/s10953-016-0447-0>

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