

QSAR model for Thyroid Receptor Alpha (TR α) antagonism (v1.0)

ProtoED

ProtoED is a computational tool designed to predict whether a compound will act as an agonist or antagonist on various hormonal receptors, facilitating the assessment of the compound's potential to disrupt the endocrine system.

By employing QSAR models, ProtoED offers an efficient alternative to experimental assays by enabling rapid and accurate predictions of compound-receptor interactions, serving as a valuable tool in chemical and pharmacological research.

This module promotes the use of alternative methods, helping to reduce the need for *in vivo* testing and supporting decision-making processes regarding potential risks to human health and the environment.

Endpoint

Human health effects: Thyroid receptor alpha antagonism

Thyroid hormone receptor alpha (TR α) is defined as a type of receptor that binds to the active thyroid hormone T3, primarily expressed in certain brain areas and the heart. Thyroid receptor alpha antagonism occurs when compounds bind to TR α but do not activate it, effectively blocking thyroid hormone signaling and potentially leading to metabolic and developmental disturbances.

Metrics

Training set

| Experimental values | QSAR predictions | |
|---------------------|------------------|------------|
| | inactive | antagonist |
| inactive | 188 | 28 |
| antagonist | 7 | 203 |

Validation set


| Experimental values | QSAR predictions | |
|---------------------|------------------|------------|
| | inactive | antagonist |
| inactive | 61 | 13 |
| antagonist | 3 | 67 |

| Parameters | Training | Validation |
|----------------------------------|----------|------------|
| Accuracy | 0.92 | 0.89 |
| Sensitivity / recall | 0.97 | 0.96 |
| Specificity | 0.87 | 0.82 |
| Precision | 0.88 | 0.84 |
| Negative predictive value | 0.96 | 0.95 |
| F-score | 0.92 | 0.89 |
| Matthews Correlation Coefficient | 0.84 | 0.79 |
| Critical Success Index | 0.85 | 0.81 |
| Area under the ROC | 0.92 | 0.89 |

ProtoED is part of



ProtoPRED platform allows the easy, fast and user-friendly prediction of different properties of chemical compounds, using proprietary (Q)SAR models.

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