In silico toxicology predictions for regulatory purposes: introduction to (Q)SAR and read-across

ProtoQSAR

Monday, October 23

	Module 1: Introduction to computational methods in regulatory chemistry
15:00-16:00	1.1 Overview of computational methods
16:00-16:30	1.2 Application of *in silico* methods in regulations (REACH, CLP, BPR, etc)
16:30-17:00	Break
	Module 2: Analogs identification and read-across with QSAR Toolbox
17:00-18:00	2.1 Analogs identification and categorization
18:00-19:00	2.2 Read-across

Tuesday, October 24

	Module 3: QSAR with different in silico platforms
15:00-16:30	3.1 Overview of QSAR predictions tools: OPERA, VEGA,EPISUITE, etc.
16:30-17:00	Break
17:00-18:00	3.2 QSAR predictions in QSAR Toolbox
18:00-19:00	3.3 QSAR predictions with ProtoPRED for regulatory purposes