



Selection guide

# Biacore™ systems

Biacore systems are designed to provide valuable information and high-quality interaction data of ions, small molecules, multidomain proteins, or viruses with targets. The systems are used in a range of scientific fields including: basic biological research, drug discovery and development, immunogenicity studies, vaccine development, and quality control. Biacore systems enable you to:

- Understand the relationship between molecular interaction and function or to confirm your results from other techniques
- Screen and characterize for hits and optimize leads based on selectivity, affinity, and kinetics
- Screen and characterize antibodies and proteins based on yes/no binding, affinity, and kinetics from the fastest on-rates to the slowest off-rates
- Quantitate protein by measuring the concentration of active protein with retained biological function



# Which system is right for you?



Applications	Biacore 8K+	Biacore 8K	Biacore S200
Kinetics/affinity characterization	●	●	●
Kinetics/affinity screening	●	●	●
Single cycle kinetics	●	●	●
LMW interaction analysis	●	●	●
Fragment screening	●	●	●
Epitope mapping	○	○	○
Immunogenicity	-	-	-
Concentration analysis	●	●	-
Parallel line analysis	●	●	-
Calibration-free concentration analysis	-	-	-
Thermodynamics	-	-	●
Comparability	-	-	-
Sample recovery MS	-	-	-
Built-in knowledge base	-	-	-

  

Specifications*	Biacore 8K+	Biacore 8K	Biacore S200
Association rate ( $k_a$ )	Proteins: up to $10^9 \text{ M}^{-1}\text{s}^{-1}$ LMW molecules: up to $10^7 \text{ M}^{-1}\text{s}^{-1}$		Proteins: $10^3$ to $3 \times 10^9 \text{ M}^{-1}\text{s}^{-1}$ LMW molecules: $10^3$ to $5 \times 10^7 \text{ M}^{-1}\text{s}^{-1}$
Dissociation rate ( $k_d$ )	$10^{-6}$ to $0.5 \text{ s}^{-1}$		$10^{-5}$ to $2 \text{ s}^{-1}$
Affinity range	fM to mM		fM to mM
Concentration range	LOD < 1 pM		-
Precision (concentration analysis)	< 5% CV		-
Molecular weight limit	No lower limit		No lower limit
Baseline noise	Typically < 0.02 RU (RMS)		< 0.015 RU (RMS)
Baseline drift	Typically < 0.3 RU/min		< 0.3 RU/min
Sample volume (kinetics)	Injection volume plus 20 to 50 $\mu\text{L}$ (application dependent)		50 to 400 $\mu\text{L}$
Immobilized molecule consumption	Typically 0.03 to 3 $\mu\text{g}/\text{flow cell}$		0.03 to 3 $\mu\text{g}/\text{flow cell}$
Analysis temperature	4°C to 40°C		4°C to 45°C
Sample storage temperature	4°C to 40°C		4°C to 45°C
Data collection rate	1 or 10 Hz		1, 10, or 40Hz
Sample capacity	12 x 96- or 384-well microplates, normal and deep-well	4 x 96- or 384-well microplates, normal and deep-well	1 x 96 or 384
Number of flow cells	16 in 8 channels		4
Unattended run time	72 h	60 h	48 h
Data evaluation time	96 curve sets < 5 min		32 curve sets < 5 min
GxP support	Yes (GxP Extension software package)		-
Additional packages	Biacore Insight Evaluation Software and application specific packages • Extended Screening and Characterization • Concentration and Potency • GxP Extension		Biacore Insight Evaluation Software and application specific packages • Extended Screening and Characterization • Concentration and Potency

● Application may be performed using the indicated product.

○ Application may be performed, but with limitations in software and/or hardware functionality.

\* Specifications are representative values, which can vary dependent on experimental conditions and individual properties of ligand and analyte.



### Biacore T200

### Biacore X100

### Biacore C

●	●	-
●	●	-
●	●	-
●	○	-
●	-	-
○	○	-
●	○	-
●	●	●
○	-	-
●	●	-
●	○	-
●	-	-
●	-	-
-	●	-

Proteins: $10^3$ to $3 \times 10^9$ M <sup>-1</sup> s <sup>-1</sup> LMW molecules: $10^3$ to $5 \times 10^7$ M <sup>-1</sup> s <sup>-1</sup>	Proteins: $10^3$ to $10^7$ M <sup>-1</sup> s <sup>-1</sup>	N/A
$10^{-5}$ to $1$ s <sup>-1</sup>	$10^{-5}$ to $0.1$ s <sup>-1</sup>	N/A
fM to mM	fM to mM	N/A
1 pM to 2 mM	10 pM to 2 mM	10 pM to 1 mM
< 5% CV	< 5% CV	< 5% CV
No lower limit	M <sub>r</sub> > 100	M <sub>r</sub> > 180
< 0.03 RU (RMS)	< 0.1 RU (RMS)	< 0.6 RU (RMS)
< 0.3 RU/min	< 0.3 RU/min	< 0.3 RU/min
50 to 400 μL	32 to 120 μL	5 to 325 μL (concentration analysis)
0.03 to 3 μg/flow cell	Typically 1 μg	Typically 1 μg
4°C to 45°C	Ambient/4°C to 40°C (Plus)	25°C (fixed)
4°C to 45°C	Ambient	Ambient (external thermostating possibility)
1 or 10 Hz	1 Hz	1 Hz
1 × 96 or 384	15 vials	2 × 96
4	2	4
48 h	24 h	24 h
32 curve sets < 10 min	1 curve set < 0.5 min	Total analysis time < 5 min/sample
Yes (Biacore T200 GxP Software Package)	-	Yes (Included)
Biacore Insight Evaluation Software and application-specific packages • Extended Screening and Characterization • Concentration and Potency • GxP Extension	Plus	-

# Label-free interaction analysis —from research, through drug discovery and development, to QC

## **Biacore 8K+**

**Discover more, with maximized capacity**



Efficiently delivers binding data of outstanding quality, meeting your toughest challenges in screening, characterization, process optimization, and quality control.

- Single solution for interaction analysis allowing ranking, kinetics, affinity, concentration, and relative potency
- Screening of 2300 molecules in a day
- High-quality kinetic characterization of 64 interactions in 4 h
- Exceptional sample capacity of up to 12 × 384-well microplates
- 72 h unattended run time with queueing abilities and rapid multirun evaluation

## **Biacore 8K**

**Discover more, more efficiently**



Efficiently delivers binding data of outstanding quality, meeting your toughest challenges in screening, characterization, process optimization, and quality control.

- Single solution for interaction analysis allowing ranking, kinetics, affinity, concentration, and relative potency
- Screening of 2300 molecules in a day
- High-quality kinetic characterization of 64 interactions in 4 h
- High sample capacity of up to 4 × 384-well microplates
- 60 h unattended run time with queueing abilities and rapid multirun evaluations

## **Biacore S200**

**Empowers low-molecular weight drug discovery programs**



Our most sensitive SPR-based sensor for cutting-edge applications.

- Determine affinity and kinetics with exceptional quality for confident lead optimization even for difficult targets where response levels are low
- Find site-selective binders directly using competition experiments
- Obtain fragment binding data from 384 single-concentration samples in less than 16 h

## **Biacore T200**

**Versatility all the way from research to discovery and quality control**



Versatile system for high-quality characterization of molecular interactions—from ions to viruses.

- Increase understanding of molecular mechanisms and structure-function relationships
- Select and optimize lead compounds during drug discovery
- Select, characterize, and assess comparability of biotherapeutics
- Perform time- and cost-efficient concentration analysis

## **Biacore X100**

**Boost your protein interaction research**



Reliable insights into biological processes in multiuser environments/small-scale interaction analysis.

- Characterize molecular mechanisms and interaction pathways based on kinetics and affinity
- Gain increased understanding of structure-function relationships
- Determine the active concentration without the need for standard curves

## **Biacore C**

**Confident concentration analysis in GxP environments—from preclinical development to QC**



The world's first SPR instrument dedicated to rapid determination of concentration in GxP environments.

- Determine the active concentration of mAbs, protein therapeutics, and vaccines or quantitate process-specific impurities
- Monitor quality and stability in formulations development and protein manufacturing
- Perform biopharmaceutical release testing

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