More Chemistries, More Choices For Solving Your Toughest Separation Challenges

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The InfinityLab Poroshell 120 family has grown to include 3 particle sizes and 20 chemistries, so you can efficiently separate the widest variety of compounds.







| InfinityLab Poroshell 120 | Chemistry | Particle Sizes | Pore Size | Temperatur Limit | e pH Range | Endcapped | Carbon Load | Surface Area | USP Designation | Benefits and Applications |
|------------------------------|-------------------------------------------------------------------------------------------------------------|----------------------|--------------|---------------------|---------------|-------------|----------------|-----------------|--------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| EC-C18 | O — Si CH ₃ CH ₃ | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 10% | 130 m2/g | L1 | General purpose Excellent peak shape and efficiency for acids, bases, and neutrals |
| EC-C8 | O — Si ~ CH ₃ CH ₃ | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 5% | 130 m2/g | L7 | General purpose Lower retention of hydrophobic analytes vs. C18 |
| Aq-C18 | -0—Si | 2.7 μm | 120 Å | 90 °C | 1.0-8.0 | Yes | Proprietary | 130 m2/g | L1 | Enhanced retention for challenging polar compounds while also separating non-polar analytes 100% aqueous mobile phase compatibility and low pH stability |
| SB-C18 | R O—Si R | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 90 °C | 1.0-8.0 | No | 9% | 130 m2/g | L1 | Excellent stability at low pH Great peak shape in highly acidic conditions |
| SB-C8 | R O—Si | 2.7 μm | 120 Å | 80 °C | 1.0-8.0 | No | 5.5% | 130 m2/g | L7 | Excellent stability at low pH Lower retention of hydrophobic analytes vs. C18 |
| HPH-C18 | - 0 - Si CH ₃ CH ₃ | 1.9 μm, 2.7 μm, 4 μm | 100 Å | 60 °C | 2.0-11.0 | Yes | Proprietary | 95 m2/g | L1 | High pH capability designed for longest lifetime, especially under high pH conditions Robust performance and long lifetimes Similar selectivity compared to EC-C18 |
| HPH-C8 | CH ₃ Si CH ₃ | 2.7 μm, 4 μm | 100 Å | 60 °C | 2.0-11.0 | Yes | Proprietary | 95 m2/g | L7 | High pH capability Robust performance and long lifetimes Lower retention of hydrophobic analytes vs. C18 |
| CS-C18 | + R + R + R | 2.7 μm | 100 Å | 90 °C | 1.0-11.0 | Yes | Proprietary | 95 m2/g | L1 | High pH capability with alternate selectivity Improved peak shape and sample capacity for basic compounds with low ionic strength mobile phases |
| Bonus-RP | Polar (CH ₂) _n - Polar group | √ 2.7 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 9.5% | 130 m2/g | L60 | Alternate selectivity to C18 Unique selectivity due to a polar embedded group, stable in 100% aqueous |
| PFP | $\begin{array}{c} CH_3 \\ Si \\ CH_3 \end{array} \qquad (CH_2)_n \begin{array}{c} F \\ -F \\ F \end{array}$ | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 5.1% | 130 m2/g | L43 | Alternate selectivity Excellent peak shape for polar and nonpolar analytes Unique selectivity for aromatic and halogenated compounds |
| Phenyl-Hexyl | O — Si — CH ₃ — CH ₃ | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 9% | 130 m2/g | L11 | Alternate selectivity with aromatic groups Highly nonpolar bonded phase takes advantage of pi-pi interactions |
| SB-Aq | R R R | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 80 °C | 1.0-8.0 | No | Proprietary | 130 m2/g | L96 | Alternate selectivity Excellent peak shape and retention of polar compounds using reversed-phase LC Exceptional stability under high-aqueous conditions, including 100% water |
| EC-CN | CH_3 CH_3 CH_3 CH_3 | 2.7 μm | 120 Å | 60 °C | 2.0-8.0 | Yes | 3.5% | 130 m2/g | L10 | Alternate selectivity Use in reversed-phase for alternate selectivity of polar and mid-polar compounds Use in normal phase for excellent peak shape and retention of nonpolar analytes |
| HILIC-Z | $-R_{2} - \begin{matrix} R_{3} & 0 \\ N_{1}^{+} - R_{1} - S = 0 \\ R_{4} & 0 \end{matrix}$ | 1.9 μm, 2.7 μm, 4 μm | 100 Å | 80 °C | 2.0-12.0 | No | Proprietary | 95 m2/g | L114 | Polar analytes Excellent retention of highly polar or charged compounds by HILIC Rugged performance at high pH or high temperature |
| HILIC | | 1.9 μm, 2.7 μm, 4 μm | 120 Å | 60 °C | 1.0-8.0 | No | NA | 130 m2/g | L3 | Polar analytes Excellent retention of polar compounds by HILIC |
| HILIC-OH5 | Polyhydroxyl Cyclofructan | 2.7 μm | 120 Å | 45 °C | 1.0-7.0 | Proprietary | Proprietary | 130 m2/g | L86 | Polar analytes Fructan bonded phase offers alternate selectivity to other HILIC phases |
| Chiral-V | Vancomycin | 2.7 μm | 120 Å | 45 °C | 2.5-7.0 | Proprietary | Proprietary | 130 m2/g | L88 | Chiral separations Amines, profens, and complex basic and neutral compounds Reversed-phase, polar ionic normal phase, or polar organic modes |
| Chiral-T | Teicoplanin | 2.7 μm | 120 Å | 45 °C | 2.5-7.0 | Proprietary | Proprietary | 130 m2/g | L63 | Chiral separations Beta blockers, hydroxyl acids, amino acids, profens, benzodiazepines, and hydantoins Reversed-phase, polar ionic normal phase, or polar organic modes |
| Chiral-CD | Hydroxypropylated ß-cyclodextrin | 2.7 μm | 120 Å | 45 °C | 3.0-7.0 | Proprietary | Proprietary | 130 m2/g | L45 | Chiral separations Stimulants, fungicides, and protected amino acids Reversed phase or polar organic modes |
| Chiral-CF | Isopropyl Cyclofructan (CF6) | 2.7 μm | 120 Å | 45 °C | 3.0-7.0 | Proprietary | Proprietary | 130 m2/g | NA | Chiral separations Primary amines Polar organic or normal phase modes |

1 bar = $14.5 \, PSI$

Format

Column ID

Column length

porous columns?

InfinityLab Poroshell Chemistry

InfinityLab Poroshell 120 EC-C18

InfinityLab Poroshell 120 EC-C8

InfinityLab Poroshell 120 SB-C18

InfinityLab Poroshell 120 SB-C8

InfinityLab Poroshell 120 SB-Aq

InfinityLab Poroshell 120 EC-CN

InfinityLab Poroshell 120 HILIC

InfinityLab Poroshell 120 Bonus-RP

InfinityLab Poroshell 120 Phenyl-Hexyl

Agilent InfinityLab is an optimized portfolio of LC instruments, columns, and supplies that work together seamlessly for maximum efficiency and performance—regardless of application area. More information at:

Aligned Chemistry

ZORBAX Eclipse Plus C18

ZORBAX Eclipse Plus EC-C8

ZORBAX StableBond SB-C18

ZORBAX StableBond SB-C8

ZORBAX StableBond SB-Aq

ZORBAX Eclipse XDB-CN

ZORBAX Bonus-RP

ZORBAX HILIC Plus

ZORBAX Eclipse Plus Phenyl-Hexyl

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Which particle is best for my method?

- Maximum pressure: 1300 bar

- Ideal for: Agilent 1290 Infinity II LC

4 µm: Improved HPLC performance

- Ideal for: Agilent 1220 Infinity II LC

What column ID and length should I choose?

4.6 mm for legacy methods

What if my methods were developed on fully

3.0 mm for lower solvent use than 4.6 mm

Shorter 30 to 100 mm for fastest separations

InfinityLab Poroshell chemistries are aligned with traditional ZORBAX chemistries—making

it easy to transfer your methods from fully porous to superficially porous particle columns.

Longer 150 to 250 mm for increased resolution

2.1 mm for lowest solvent use and MS applications

For more information about InfinityLab Poroshell 120 Columns, go to www.agilent.com/chem/poroshell-120