

• Retention mainly governed by hydrophobicity, can offer other attributes but to a lesser extent

- RPLC workhorse for majority of routine assays
- Universal phase for acids, bases and neutral solutes

EUERBY-MODIFIED TANAKA PLOTS OF HALO[®] REVERSED PHASE SMALL MOLECULE CHEMISTRIES

The Euerby-modified Tanaka plot is identical minus the A parameter, as the hydrophobicity parameter (B) entails this same information. The following radar plots visually compare ten of our HALO[®] columns for reversed phase separations, normalized at each axis for comparative purposes.

EUERBY-MODIFIED TANAKA PLOTS

The Tanaka test was first developed to determine differences between C18 stationary phases. Six different parameters/column properties (A,B,C,D,E,F) are determined based on the retention factors (k) of certain solute(s) and their respective selectivity factors under different isocratic and pH conditions.

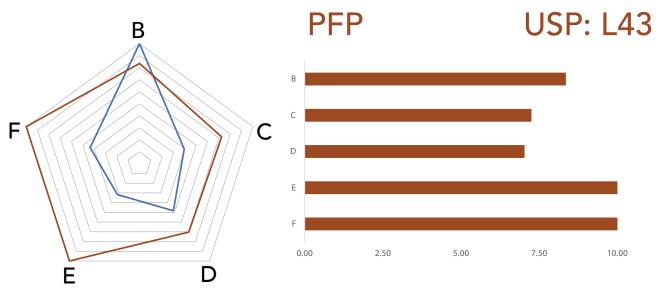
Briefly, the parameters represented are:

- A: amount of alkyl chains
- B: hydrophobicity
- C: shape selectivity
- D: hydrogen bonding capacity
- E: total ion-exchange capacity (pH>7)
- **F**: acidic ion-exchange capacity (pH<3)

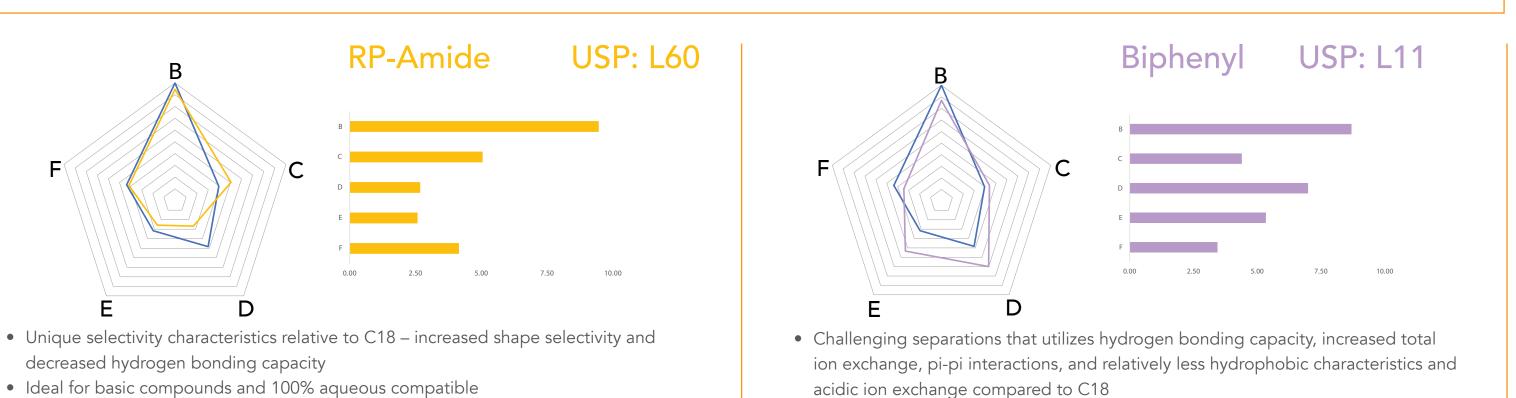
Acidic Ion-Exchange: F

Total Ion-Exchange: E

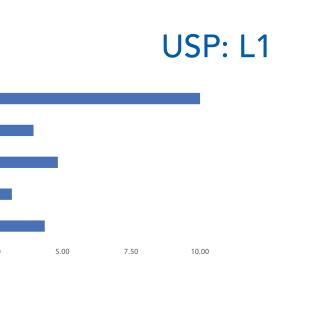
- Note: the reciprocal value was represented for elution order changes: (i) C8 – shape selectivity; (ii) Biphenyl – H- bonding capacity; and (iii) PFP – total ion-exchange.
- This global view of all the HALO[®] columns may be useful when trying to select a column based upon the five attributes. *PAH applicated phase available in ENVIROCLASS product line



- Versatility to be employed in both RP and HILIC separations and unique characteristics in comparison to C18 phase - key to include in method development/column screening work
- Highest ion exchange and total ion exchange attributes with greater shape selectivity and less hydrophobicity relative to C18
- Applications: basic drugs, mycotoxin screening, tranquilizers

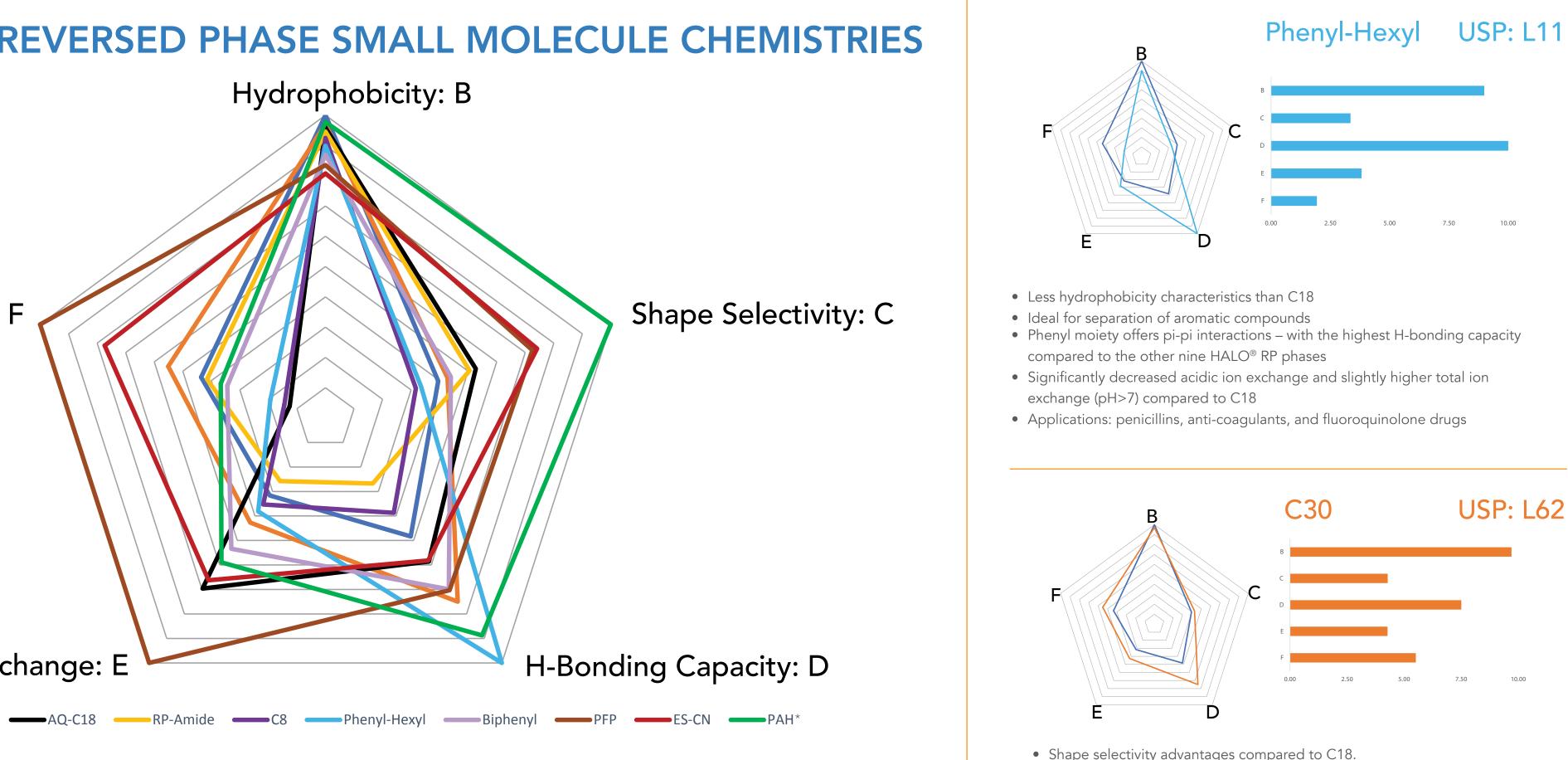


- Ideal for basic compounds and 100% aqueous compatible
- Applications: antibiotics, cholesterol lowering drugs, phenolic acids in food and beverages, active ingredients in sunscreen, melatonin and related compounds



USP: L7 **C8** C

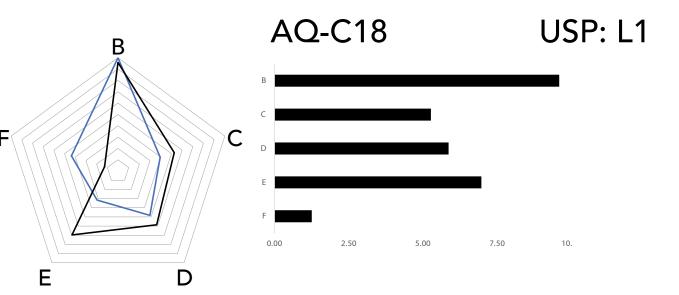
- Less hydrophobic compared to C18 and significant lower acidic ion exchange capabilities compared to C18
- Applications: cholesterol lowering pharmaceutical (statin) compounds, flavonoids, and lipid analysis of algal oil
- Ideal for broad range of analytes





• Applications: opiates, polar and non-polar pesticides and via LC-MS

- to C18



• Increased retention for polar analytes separated via RPLC and for total ion exchange capabilities

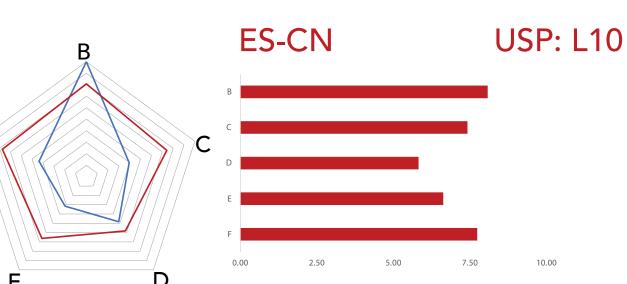
• Offers 100% aqueous compatibility

• Applications: polar pesticides, polar organic acids and separations that employ 100% water conditions e.g. pharmaceutical separation of nucleobases in 100% water 0.1%TFA <1.2 min

• Shape selectivity advantages compared to C18.

• Larger H-bonding capacity and increased ion exchange capabilities in both acidic and basic conditions compared to C18

• Applications: isomer separations, fat/water soluble vitamins, carotenoids, lipids, anti-inflammatory, anti-lymphatic and anti-allergy steroids

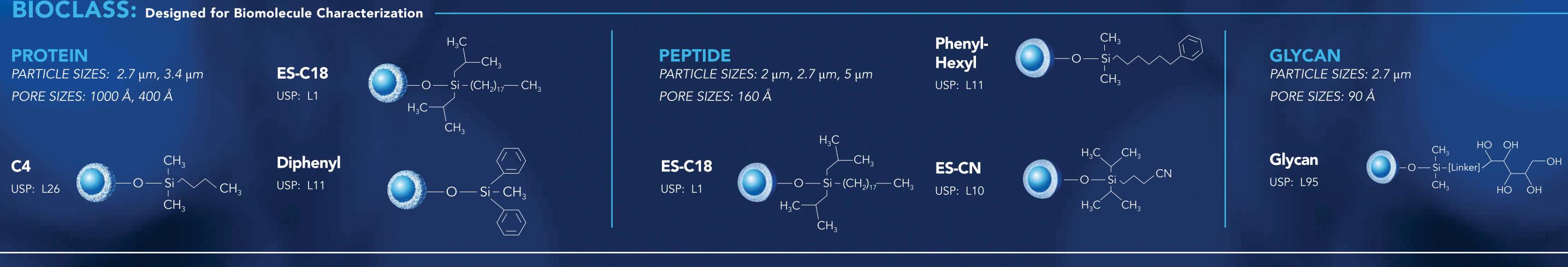


• Less hydrophobic, increased ion exchange capabilities, shape selectivity, lone pair of electrons of the phase's CN moiety provide unique RP interactions - compared

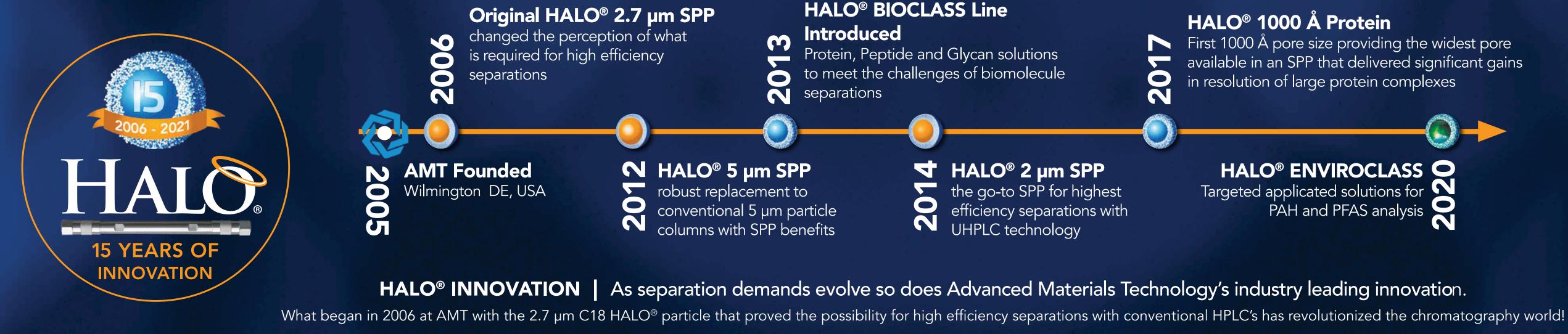
• Ideal for polar analytes and 100% aqueous compatibility making it a must have for method development/column screening

• Applications: B-lactam antibiotics, NSAIDs and penicillins





THE ORIGINAL PARTICLE WHICH FORGED A NEW PATH IN HPLC SEPARATIONS

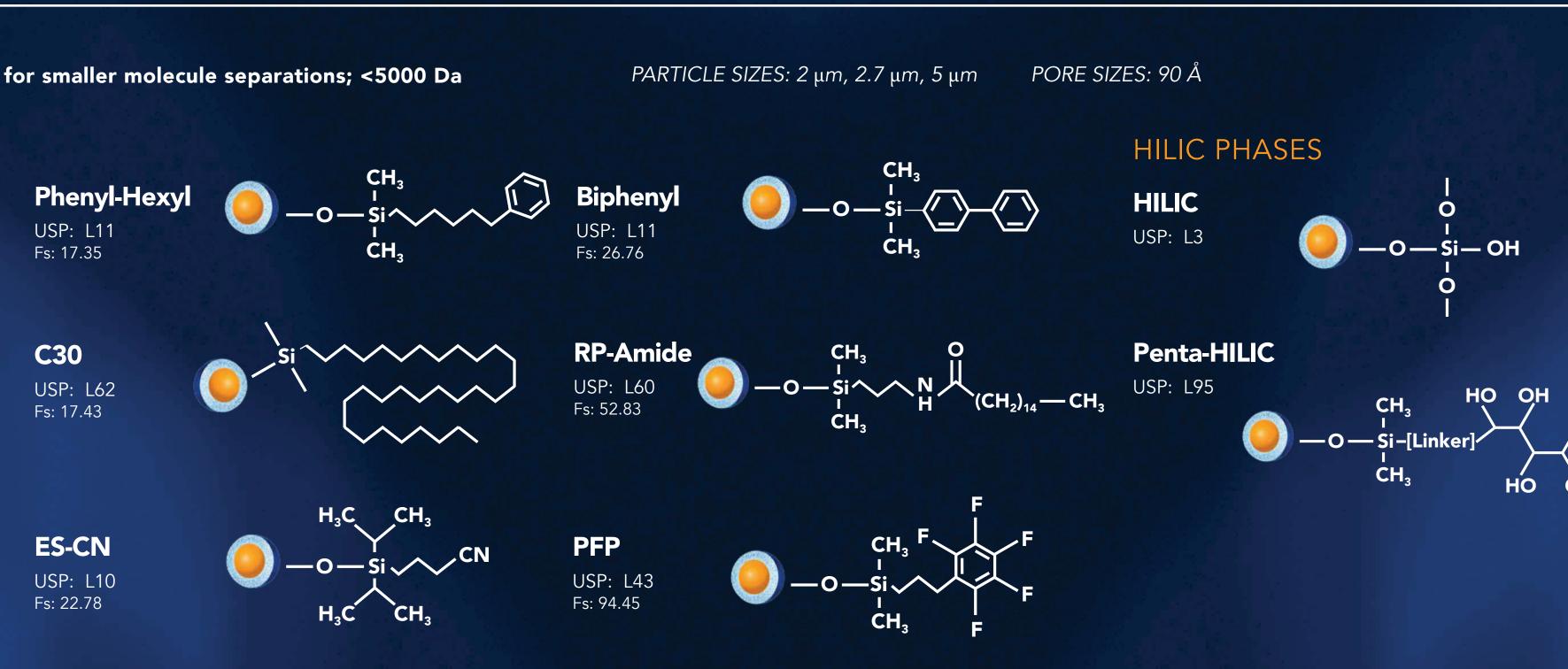


The innovation continues by forging a new path through manipulations of the particle morphology which has advanced the chromatographic separations in the biopharmaceutical and environmental industries. The timeline above represents AMT's commitment to delivering innovative tools to enhance the separation resolution afforded by the column technology.

SMALL MOLECULE: Designed for smaller molecule separations; <5000 Da where speed and resolution are critical.

REVERSED PHASE

 $-O - Si - (CH_2)_{17} - CH_3$ **C18** USP: L1 Fs: 0 CH_{3} $-O - Si - (CH_{2})_{7} - CH_{3}$ I CH_{3} **C30** USP: L62 Fs: 17.43 **C8** USP: L7 Fs: 10.04 O O Si CN PFP USP: L43 Fs: 94.45 **ES-CN** USP: L10 Fs: 22.78 AQ-C18 —[CH₂—(CH₂)₁₆—CH₃]_x [Polar Ligand], USP: L1 Fs: 12.07





INNOVATION | QUALITY | EXPERIENCE

	ENVIROLASS: Applicated solution for PFAS and PAH analysis. PARTICLE SIZE: 2.7 μm PORE SIZES: PROPRIETARY		
	PFAS Analytical	<u>()</u> -o-	CH ₃ - Si-C ₁₈ H ₃₇ - CH ₃
— ОН ОН	PFAS Delay	() -o-	CH ₃ - Si- (CH ₂) _n — CH ₃ - CH ₃
	PAH		i−(CH ₂) ₁₇ —CH ₃

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