

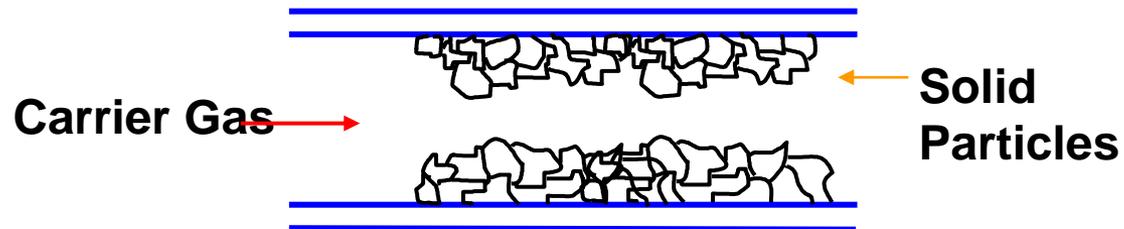
GC Method Development

Column and Phase Selection – Series 3

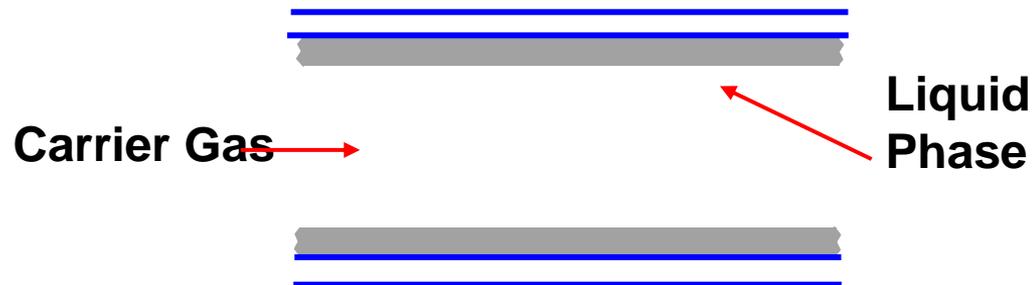
Simon Jones
Application Engineer
March 24, 2011

CAPILLARY COLUMN TYPES

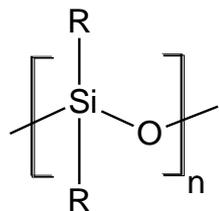
Porous Layer Open Tube (PLOT)



Wall Coated Open Tube (WCOT)

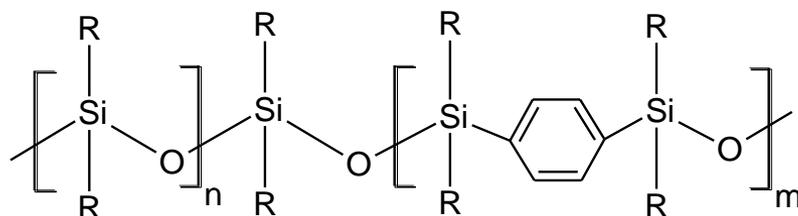


STATIONARY PHASE POLYMERS

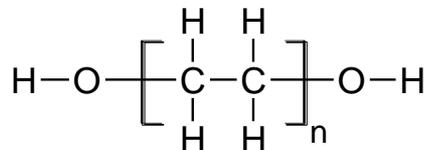


Siloxane

R=methyl, phenyl, cyanopropyl, trifluoropropyl



Siarylene backbone



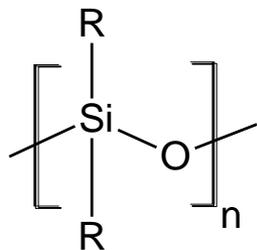
Polyethylene Glycol

Stationary Phase

% Substitution -- polysiloxanes

% = # of sites on silicon atoms occupied

Balance is methyl



Siloxane

R=methyl, phenyl, cyanopropyl, trifluoropropyl

Stationary Phase

Poly(ethylene) Glycol



100% PEG (DB-WAX)
Less stable than polysiloxanes
Unique separation characteristics

Poly(Ethylene) Glycol Modified

- Base deactivated (CAM)
- Acid Modified (DB-FFAP)
- Extended Temperature Range

Specialty Phases

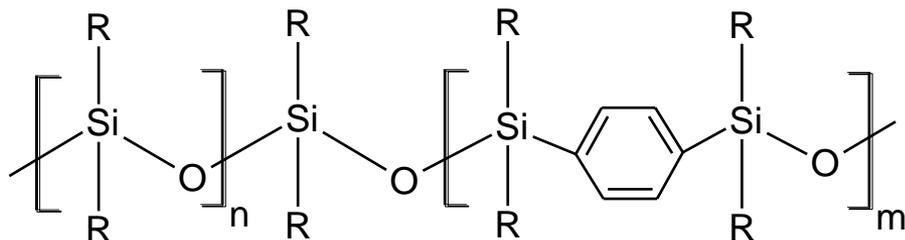
Columns developed for particular applications

Examples: DB-VRX, DB-MTBE, DB-TPH, DB-ALC1, DB-ALC2, DB-HTSimDis, DB-Dioxin, Select Low Sulfur, CP-Volamine, Select PAH, DB-EUPAH

Three Types Of Low Bleed Phases

- Phases tailored to “mimic” currently existing polymers

Examples: DB-5ms, DB-35ms, DB-17ms



Siarylene backbone

- New phases unrelated to any previously existing polymers

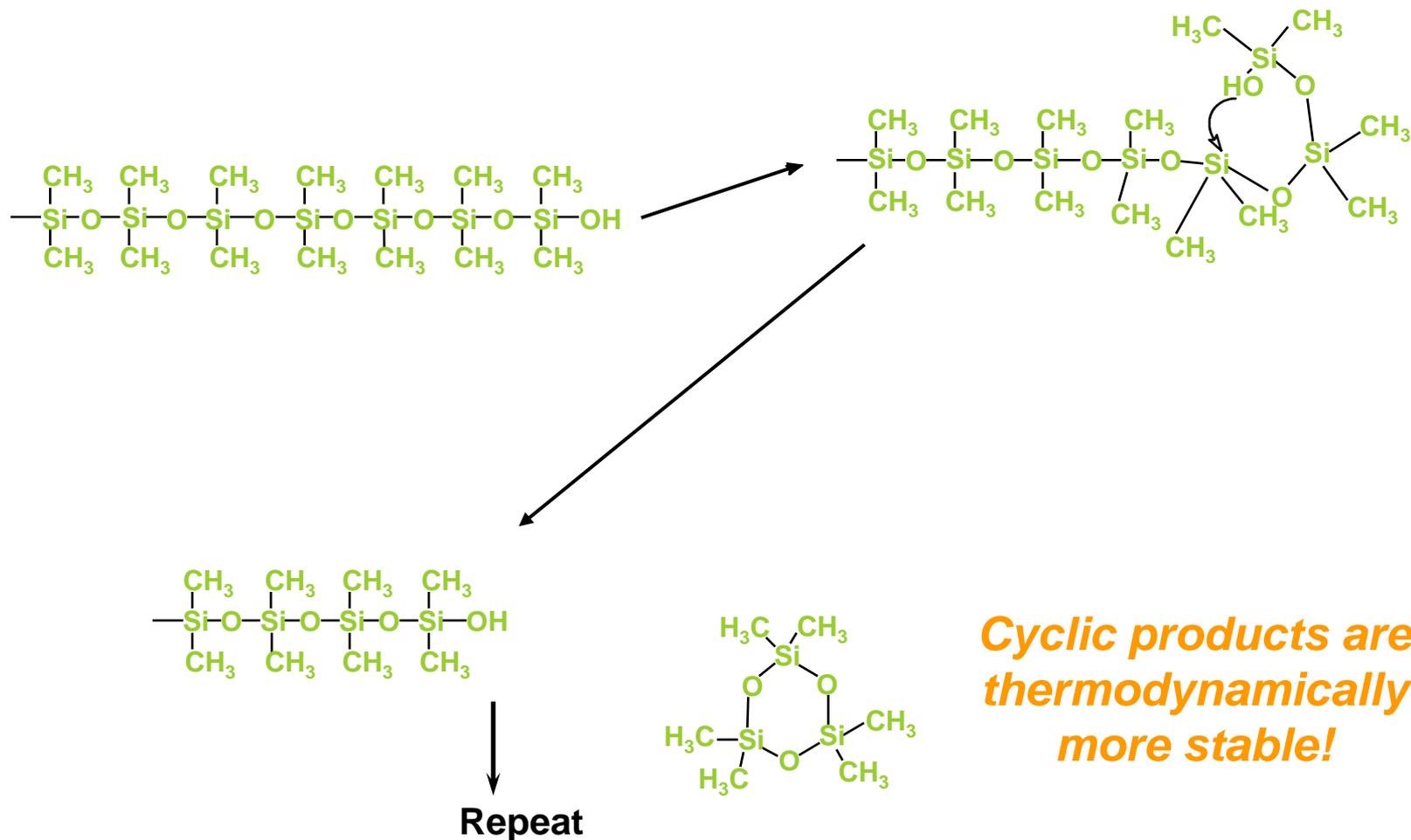
Examples: DB-XLB

- Optimized manufacturing processes

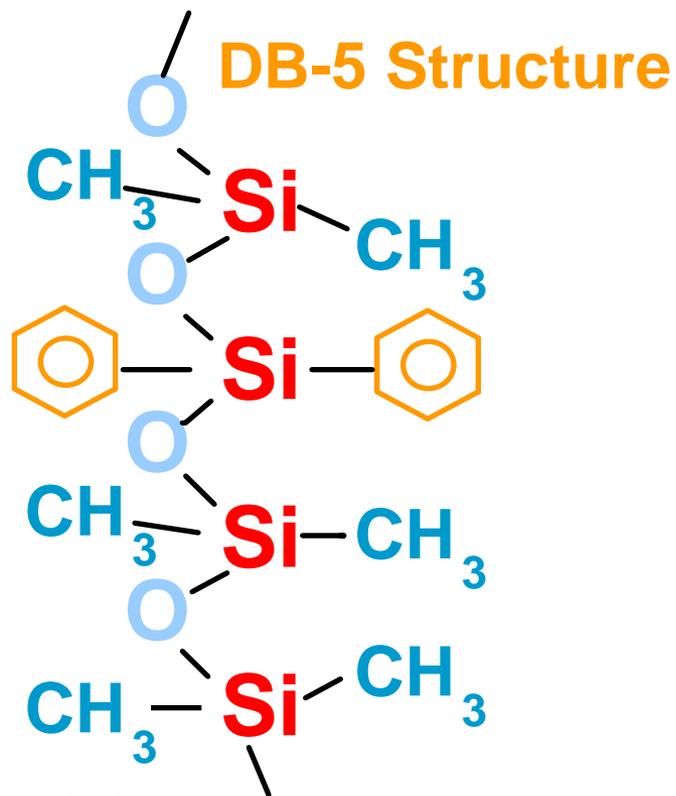
DB-1ms, HP-1ms, HP-5ms

What is Column Bleed???

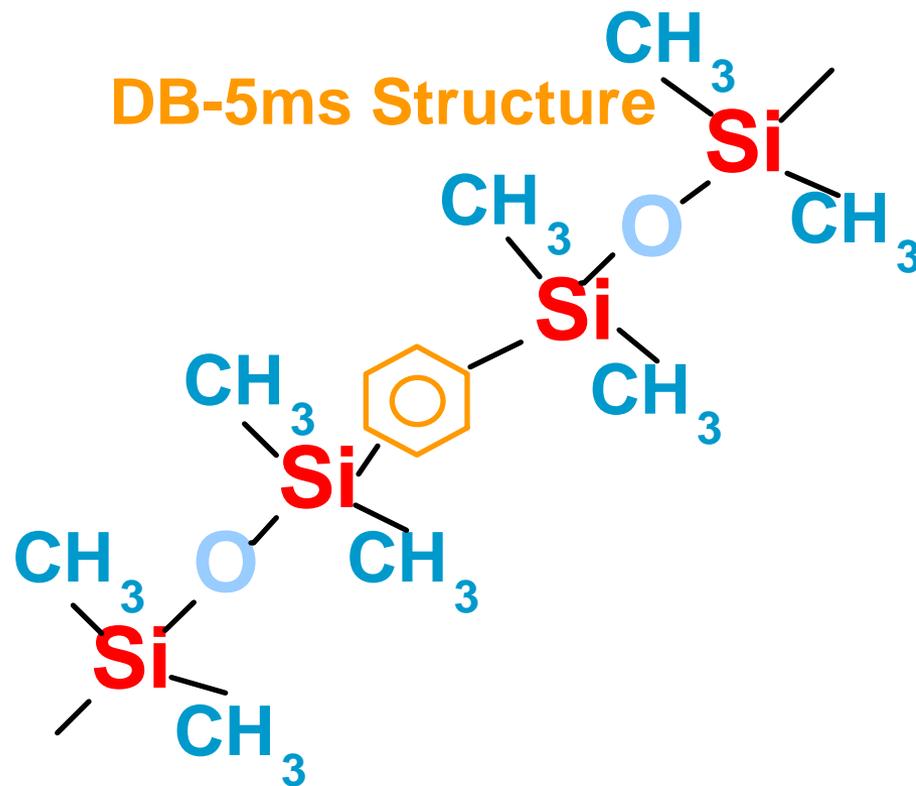
“Back Biting” Mechanism of Product Formation



DB-5ms Structure



DB-5
5% Phenyl

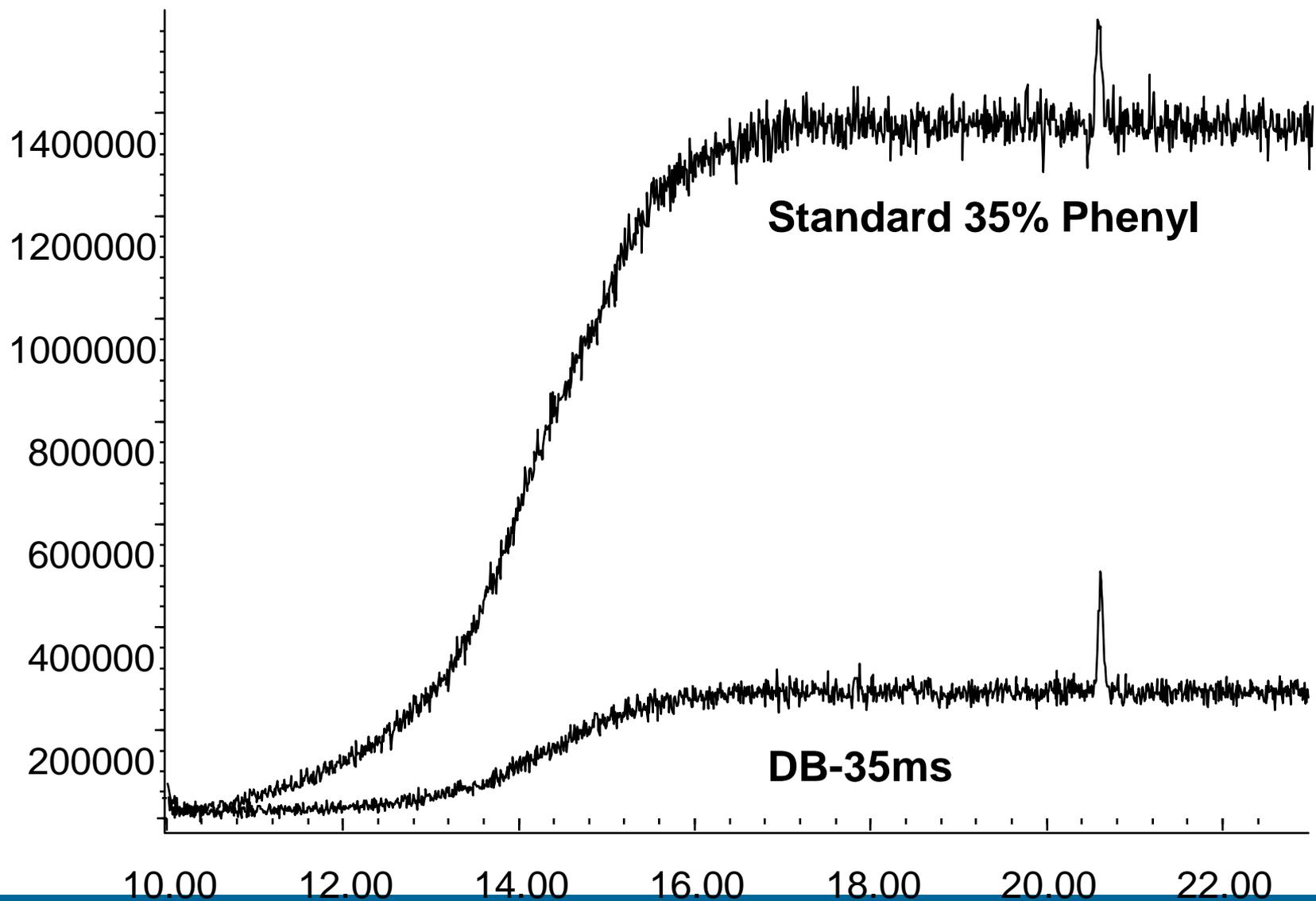


DB-5ms

1. Increased stability
2. Different selectivity
3. Optimized to match DB-5

DB-35MS VS STANDARD 35% PHENYL

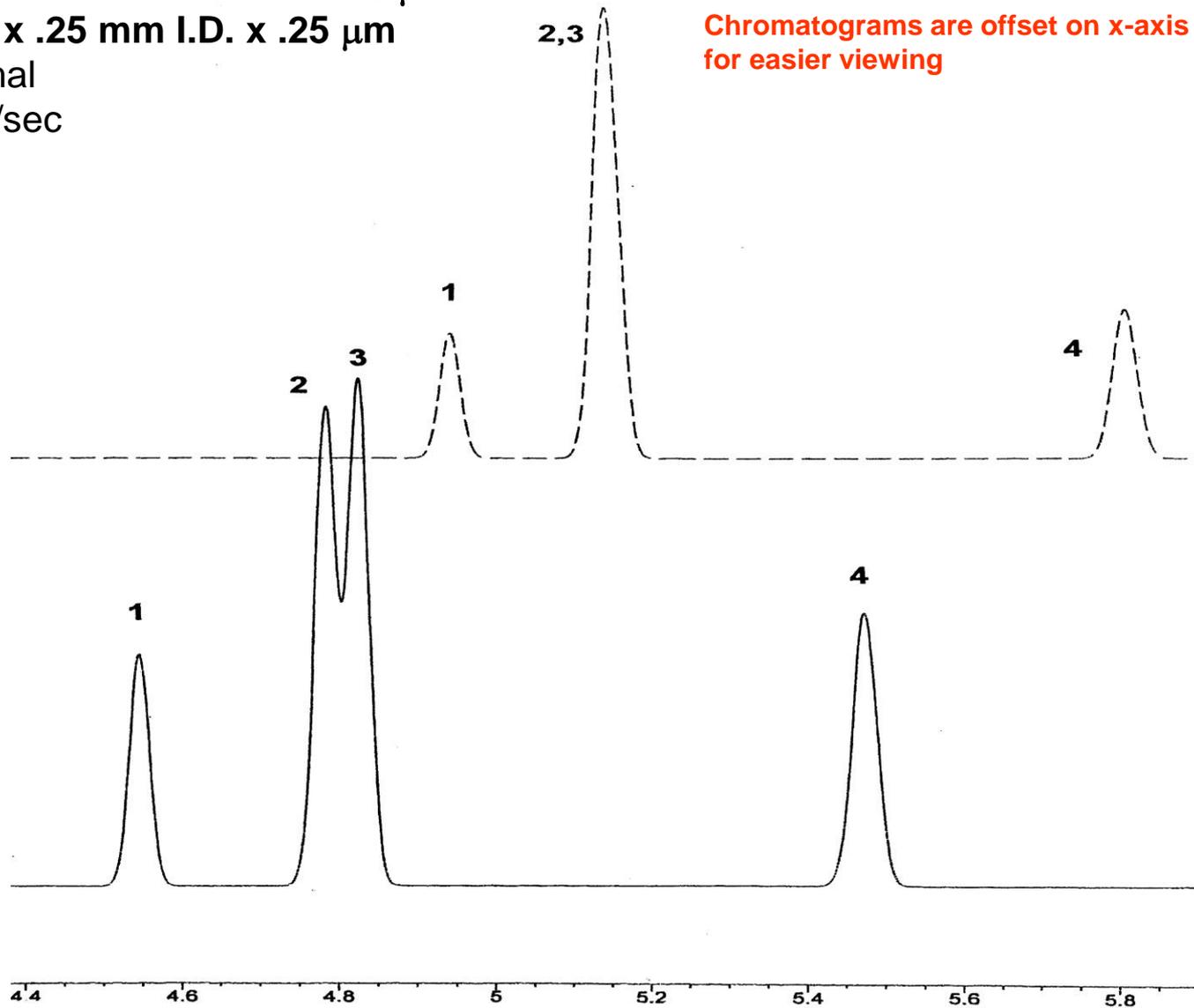
Benzo[g,h,i]perylene, 1 ng



Solid line: DB-5ms 30 m x .25 mm I.D. x .25 μ m
Dashed line: DB-5 30 m x .25 mm I.D. x .25 μ m
Oven: 60° C isothermal
Carrier gas: H₂ at 40 cm/sec

- 1: Ethylbenzene
- 2: m-Xylene
- 3: p-Xylene
- 4: o-Xylene

Chromatograms are offset on x-axis for easier viewing



Why is stationary phase type important?

Influence on α

$$\alpha = \frac{k_2}{k_1}$$

k_2 = partition ratio of 2nd peak

k_1 = partition ratio of 1st peak

Selectivity

- Relative spacing of the chromatographic peaks
- The result of all non-polar, polarizable and polar interactions that cause a stationary phase to be more or less retentive to one analyte than another

Optimizing Selectivity (α)

Match analyte polarity to stationary phase polarity

- 'like dissolves like'

Take advantage of unique interactions between analyte and stationary phase functional groups

Analyte Polarity

Nonpolar Molecules - generally composed of only carbon and hydrogen and exhibit no dipole moment (Straight-chained hydrocarbons (n-alkanes))

Polar Molecules - primarily composed of carbon and hydrogen but also contain atoms of nitrogen, oxygen, phosphorus, sulfur, or a halogen (Alcohols, amines, thiols, ketones, nitriles, organo-halides, etc. Includes dipole-dipole interactions and H-bonding)

Polarizable Molecules - primarily composed of carbon and hydrogen, but also contain unsaturated bonds (Alkenes, alkynes and aromatic compounds)

Selectivity Interactions

- Dispersion
- Dipole
- Hydrogen bonding

Dispersion Interaction

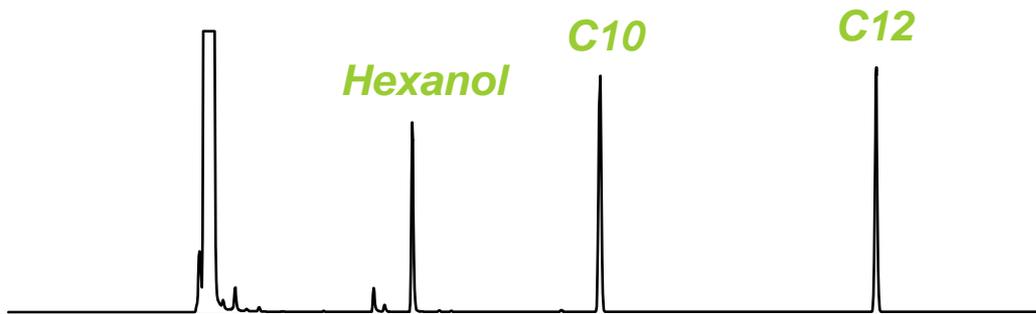
ΔH_{vap}

- Separation by differences in analyte heat of vaporizations (ΔH_{vap})
- Heat necessary to convert a liquid into a gas (at the same temperature)

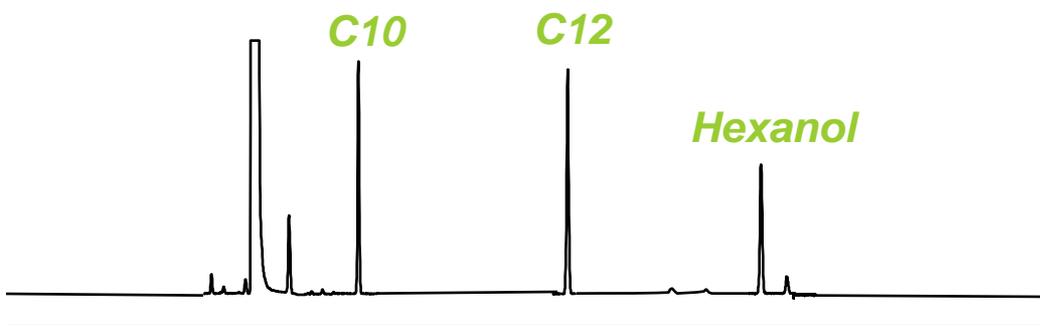
Dispersion Interaction

Solubility And Retention

Hexanol 158°C
Decane 174°C
Dodecane 216°C



**100% Methyl
(non-polar)**



**100% PEG
(polar)**

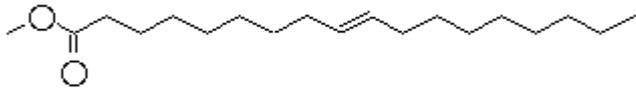
30 m x 0.32 mm ID, 0.25 μ m
He at 35 cm/sec
50-170°C at 15°/min

Dispersion Interaction

$$\Delta H_{\text{vap}}$$

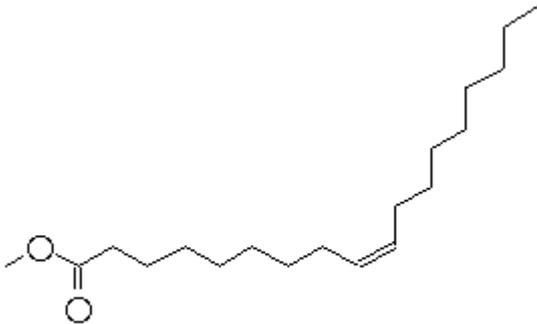
- ✓ **Vapor pressure: good approximation**
- ✓ **Boiling point: poor approximation**

Dipole Interaction



C18:1 (Methyl *trans*-9-octadecenoate)

B.Pt. 186°C



C18:1 (Methyl *cis*-9-octadecenoate)

B.Pt. 186°C

Smaller differences require a stronger dipole phase

Bacterial Fatty Acid Methyl Esters

Column: DB-23

30 m x 0.25 mm I.D., 0.25 μ m

J&W P/N: 122-2332

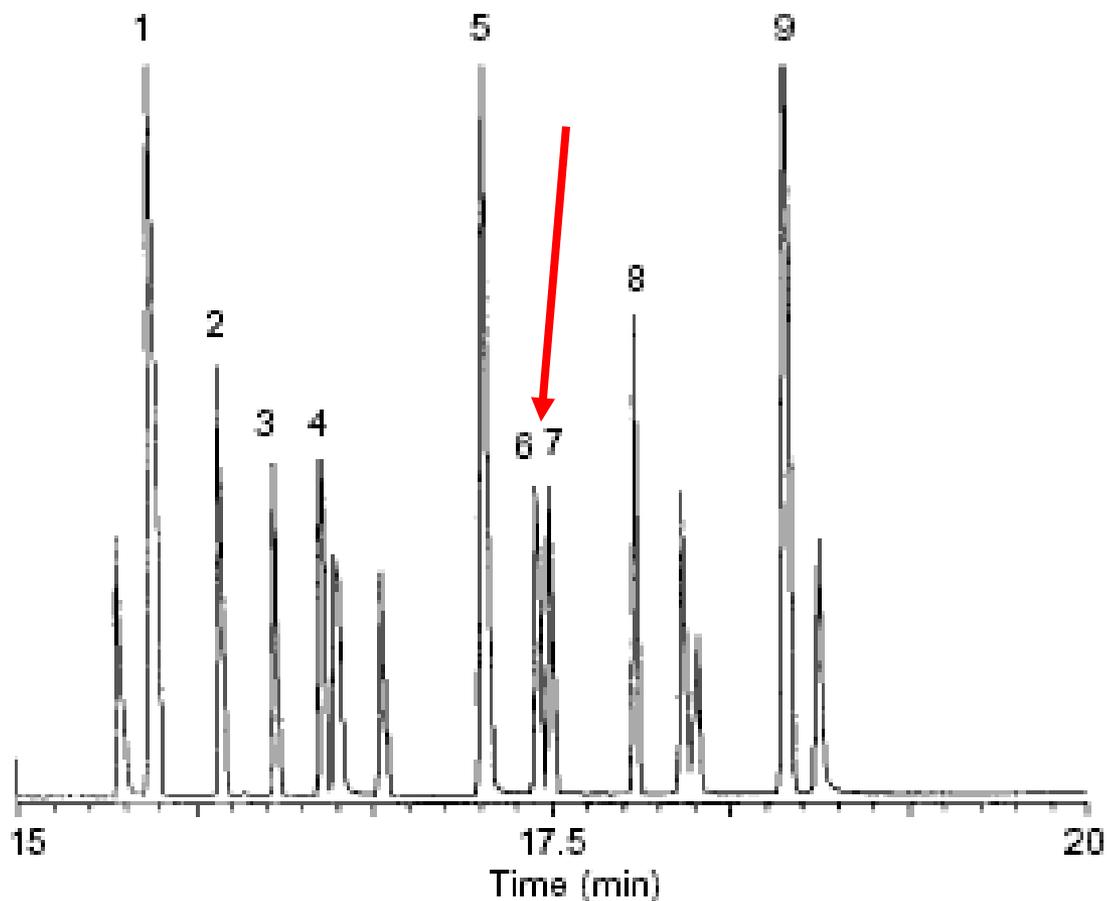
Carrier: Hydrogen at 40 cm/sec

Oven: 90°C for 6 min
90-210°C at 10°/min

Injector: Splitless, 45 sec purge activation time, 1 μ L

Detector: Finnigan INCOS 50 MSD

1.	14-Me C15:0	14-Methyl pentadecanoate
2.	C16:1	<i>cis</i> -9-Hexadecanoate
3.	14-Me C16:0	14-Methyl hexadecanoate
4.	C17:0	Heptadecanoate
5.	15-Me C17:0	15-Methyl heptadecanoate
6.	C18:1	<i>cis</i> -Octadecanoate
7.	C18:1	<i>trans</i> -Octadecanoate
8.	C19:0	Nonadecanoate
9.	C20:0	Arachidate



Fames – 37 Component Standard

Column: DB-23
60 m X 0.25 mm X 0.15 μ m

Agilent P/N 122-2361

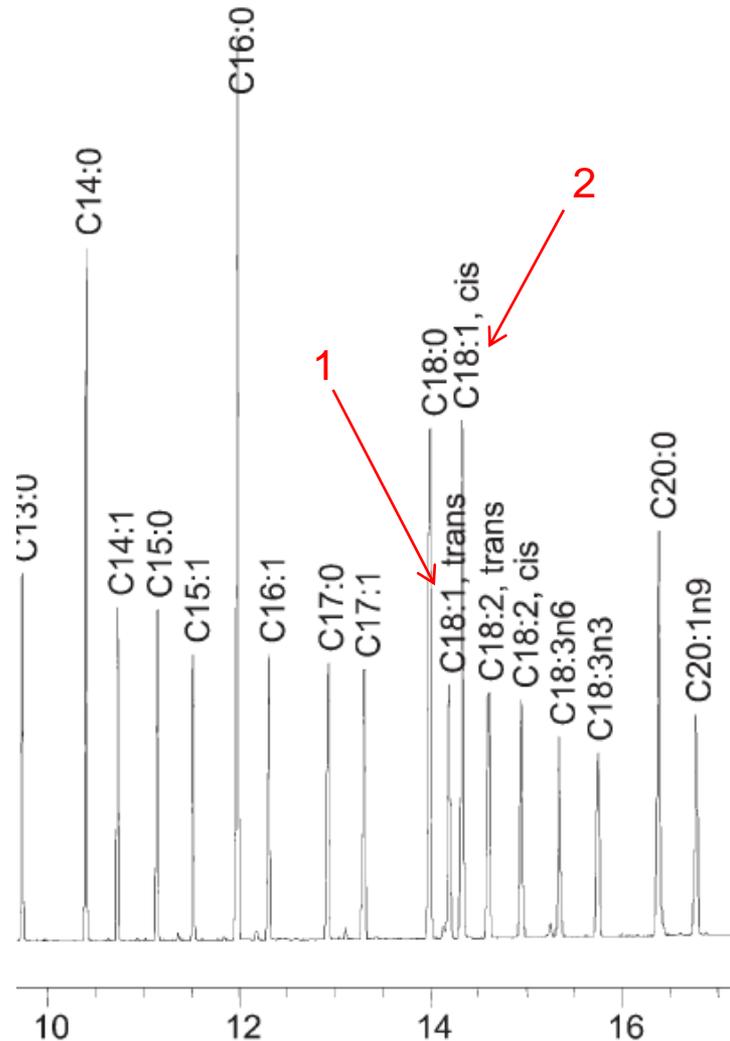
Carrier: He, 33 cm/sec @ 50°C

Oven: 50°C for 1 min
25°C/min to 175 (no hold)
4°C/min to 230°C hold 5 min

Injector: 250°C, Split 50:1, 1 μ L

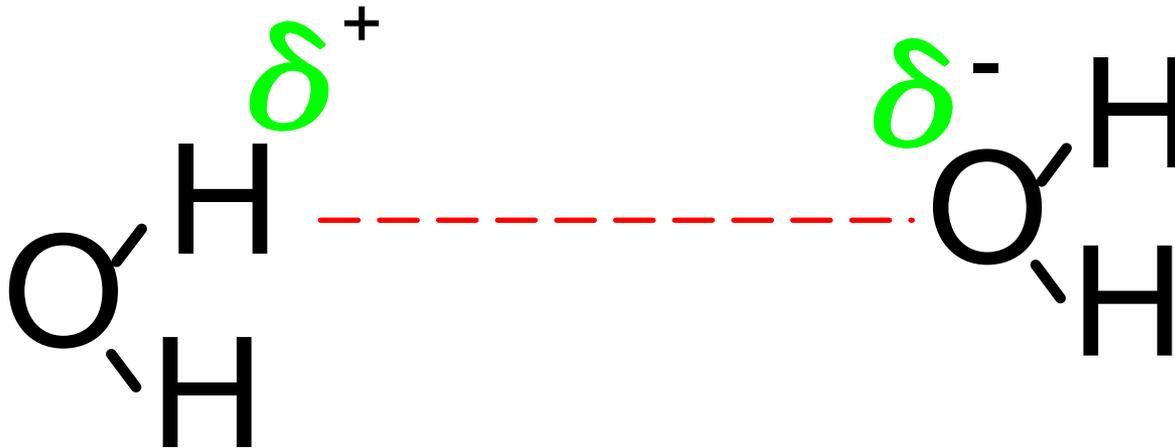
Detector: FID, 250°C

- 1 C18:1 (Methyl *trans*-9-octadecenoate)
- 2 C18:1 (Methyl *cis*-9-octadecenoate)



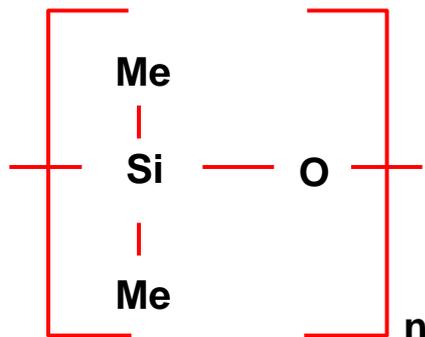
Hydrogen Bonding Interaction

Dipole-Dipole interaction with H bound to O or N interacting with an O or N



NONPOLAR PHASES

Typified by 100% polydimethylsiloxanes such as HP-1, DB-1, DB-1ms, HP-1ms, VF-1ms, CP-Sil 5 CB

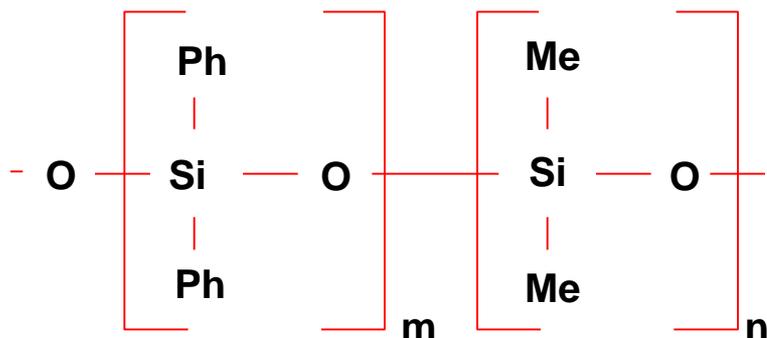


Separation Mechanisms:

- Dispersion only

POLARIZABLE PHASES

Typified by phenyl substituted siloxanes, substituted at 5-50% (HP-5, HP-5ms, DB-35, DB-35ms, DB-17, DB-17ms)



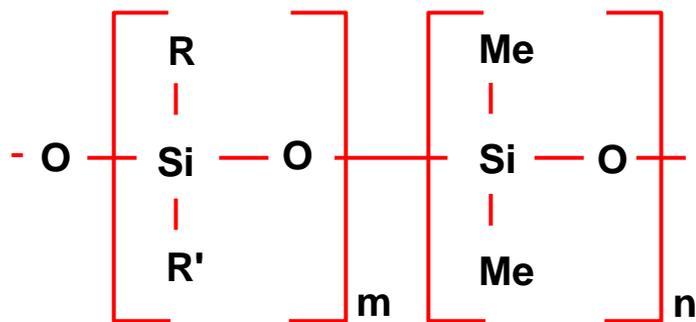
5%--weakly polar,
rest--mid polar

Separation Mechanisms:

- Dispersion
- Inducible dipole at phenyl groups

STRONG DIPOLE PHASES

Typified by cyanopropyl or trifluoropropyl substituted siloxanes, substituted 6-50% (DB-1701, DB-1301, DB-200, DB-23, DB-225)



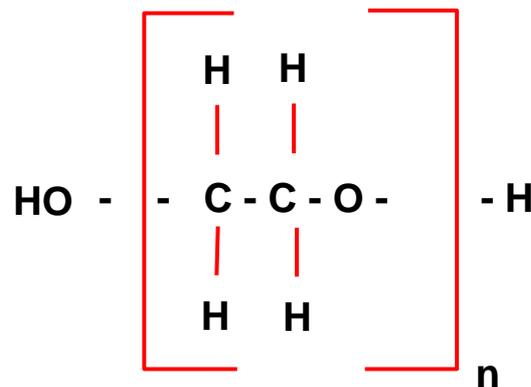
R = cyanopropyl or trifluoropropyl
R' = phenyl or methyl

Separation Mechanisms:

- Dispersion
- Inducible dipole at phenyl groups
- Strong permanent dipole
- Hydrogen bonding

HYDROGEN BONDING PHASES

Typified by polyethylene glycol polymers (Carbowax, HP-INNOWax, DB-WAX, DB-FFAP, CAM)



Separation Mechanisms:

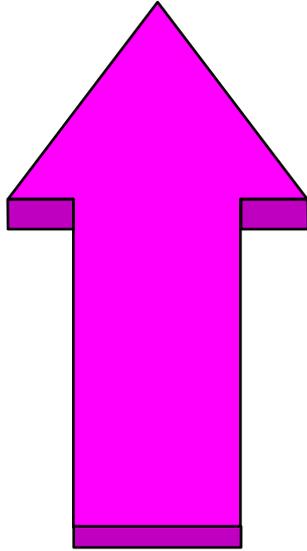
- Dispersion
- Strong permanent dipole
- Hydrogen bonding

Selectivity

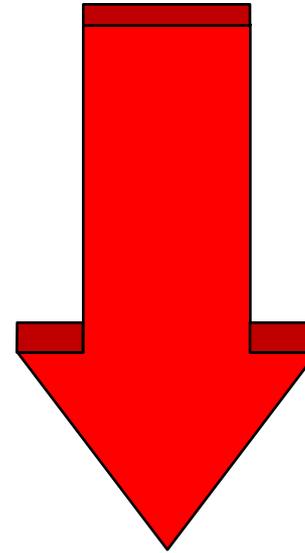
Interaction Strengths

Phase	Dispersion	Dipole	H Bonding
Methyl	Strong	None	None
Phenyl	Strong	None	Weak
Cyanopropyl	Strong	Strong	Moderate
Trifluoropropyl	Strong	Moderate	Weak
PEG	Strong	Strong	Moderate

Polarity



Polarity

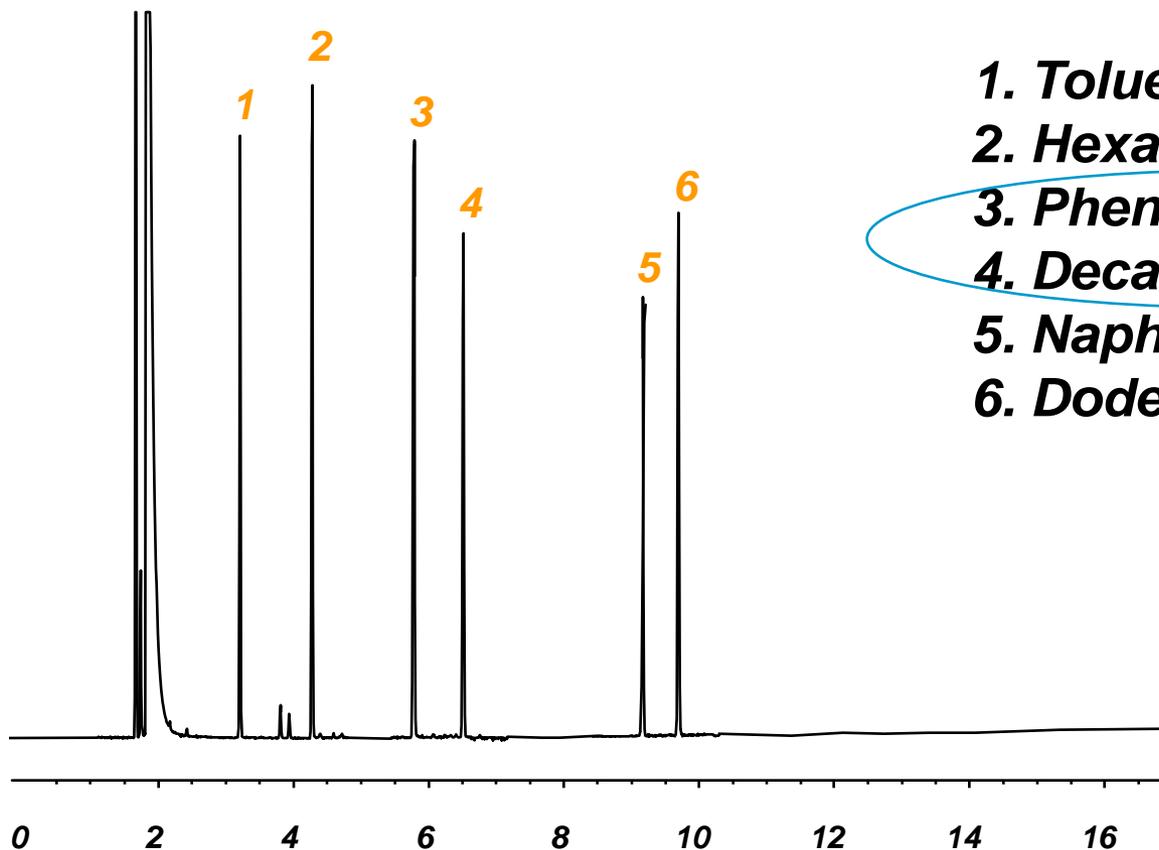


**Stability
Temperature Range**

Compounds Properties

Compounds	Polar	Aromatic	Hydrogen Bonding	Dipole
Toluene	no	yes	no	induced
Hexanol	yes	no	yes	yes
Phenol	yes	yes	yes	yes
Decane	no	no	no	no
Naphthalene	no	yes	no	induced
Dodecane	no	no	no	no

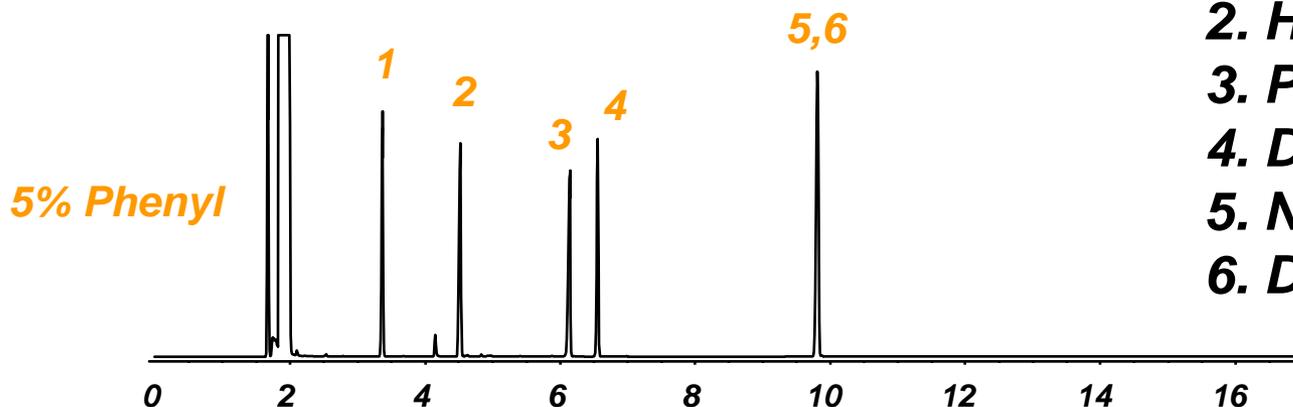
100% Methyl Polysiloxane



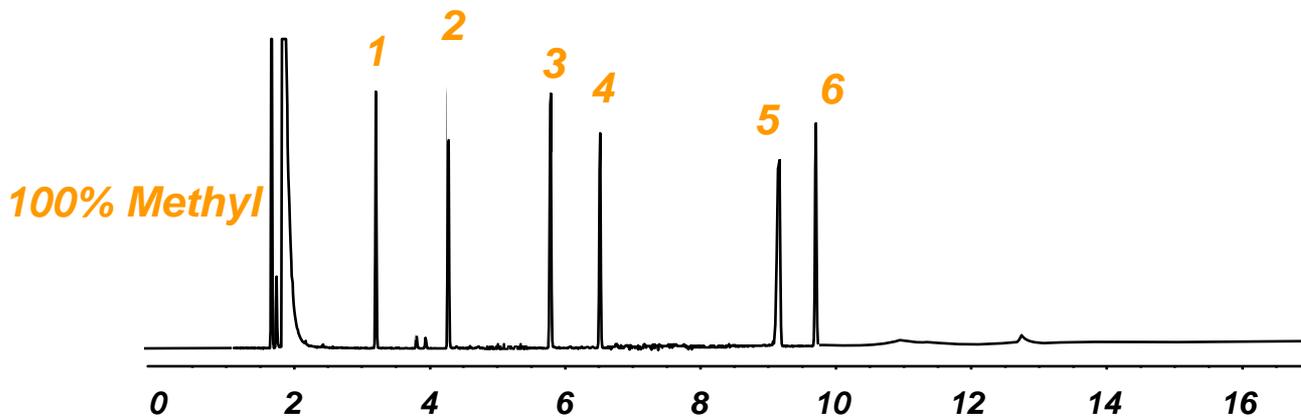
- | | |
|-------------------|-------|
| 1. Toluene | 110°C |
| 2. Hexanol | 158°C |
| 3. Phenol | 181°C |
| 4. Decane (C10) | 174°C |
| 5. Naphthalene | 218°C |
| 6. Dodecane (C12) | 216°C |

Strong Dispersion
No Dipole
No H Bonding

5% Phenyl



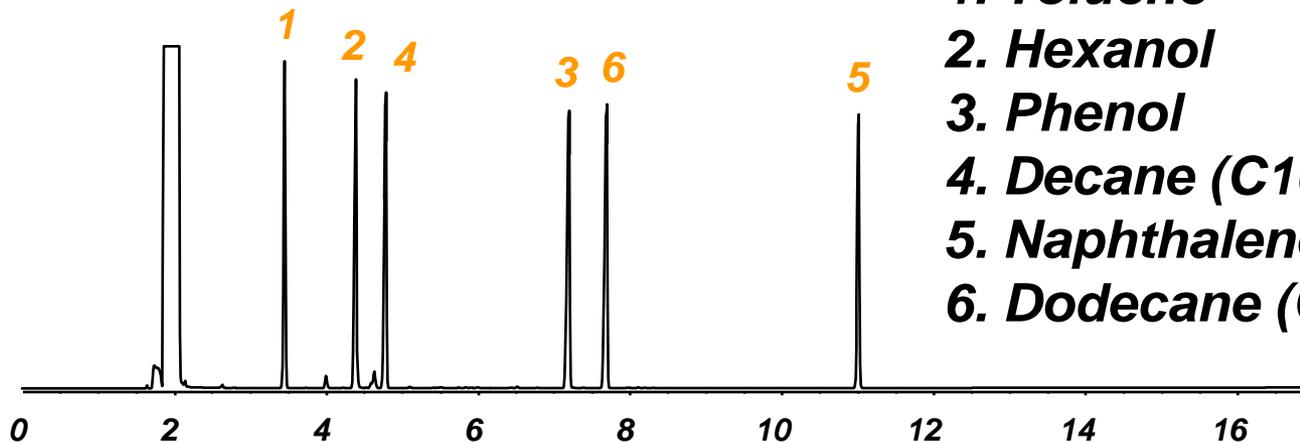
1. Toluene	110°C
2. Hexanol	158°C
3. Phenol	181°C
4. Decane (C10)	174°C
5. Naphthalene	218°C
6. Dodecane (C12)	216°C



Strong Dispersion
No Dipole
Weak H Bonding

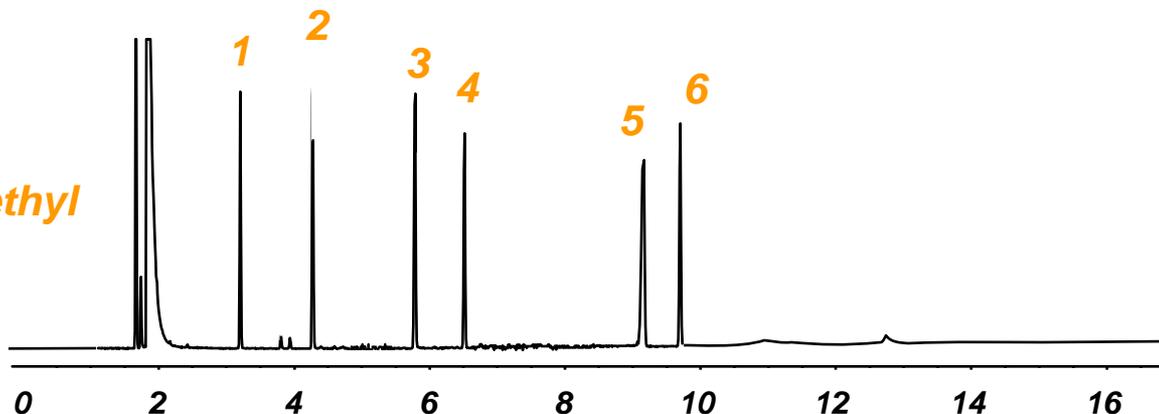
50% Phenyl

50%
Phenyl



1. Toluene 110°C
2. Hexanol 158°C
3. Phenol 181°C
4. Decane (C10) 174°C
5. Naphthalene 218°C
6. Dodecane (C12) 216°C

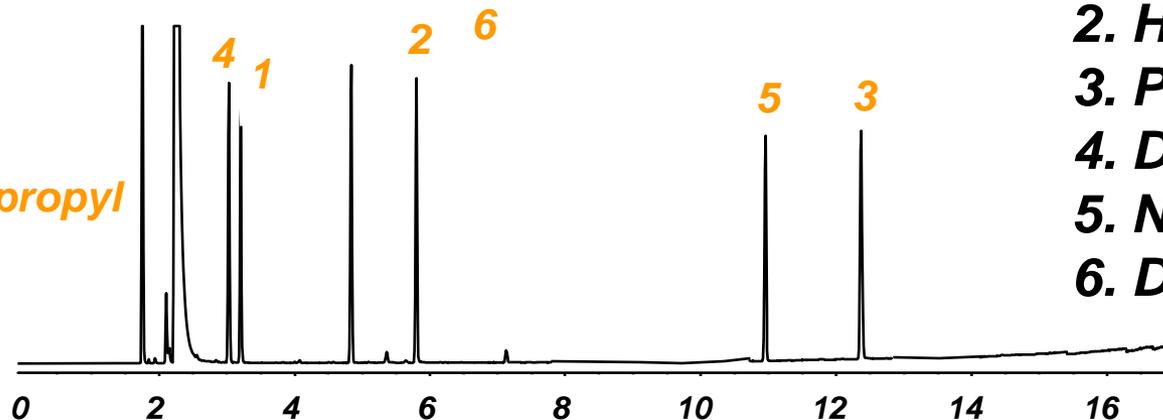
100% Methyl



Strong Dispersion
No Dipole
Weak H Bonding

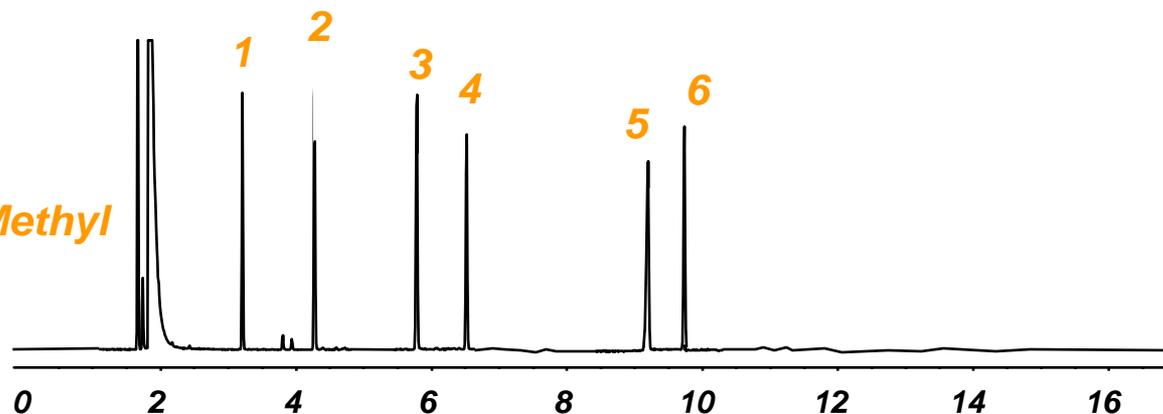
50% Cyanopropyl

50%
Cyanopropyl



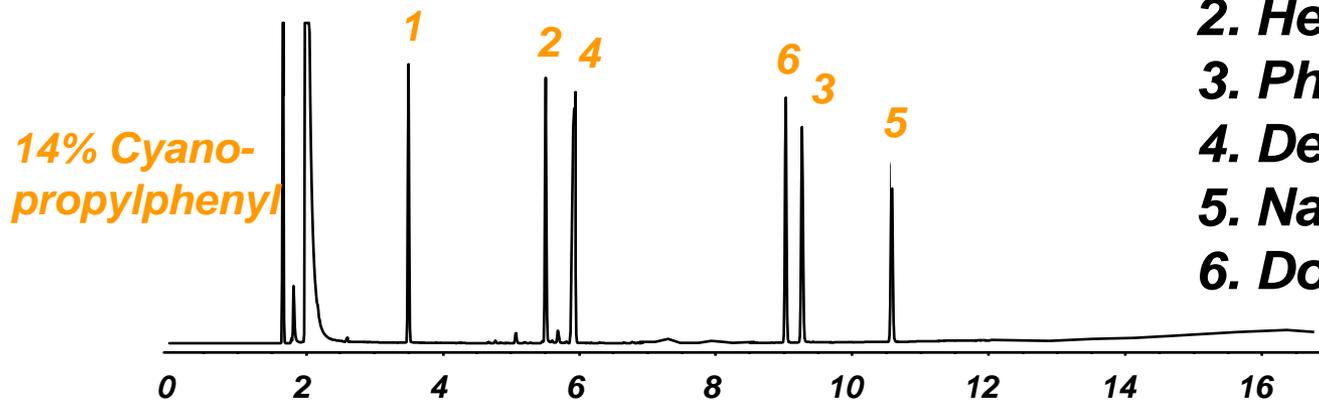
1. Toluene	110°C
2. Hexanol	158°C
3. Phenol	181°C
4. Decane (C10)	174°C
5. Naphthalene	218°C
6. Dodecane (C12)	216°C

100% Methyl

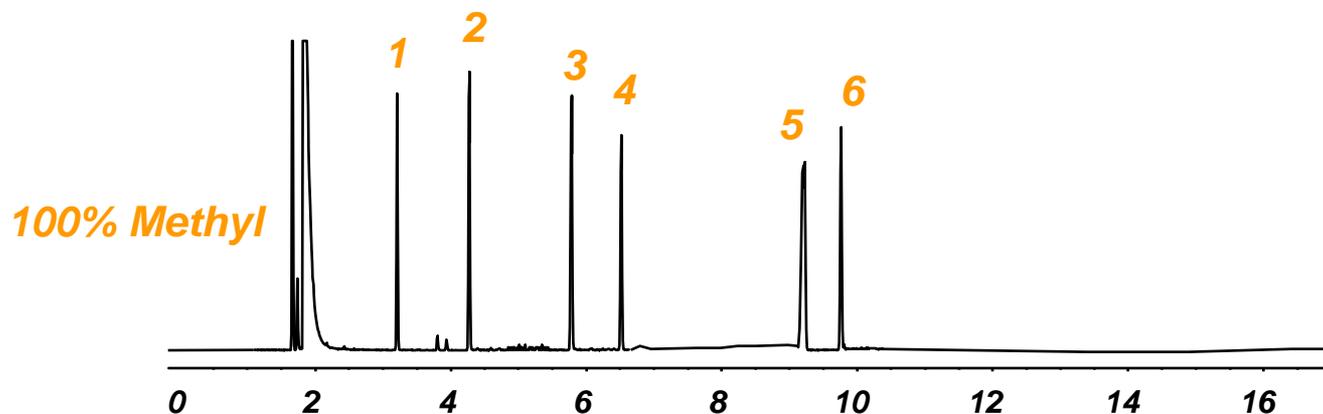


Strong Dispersion
Strong Dipole
Moderate H Bonding

14% Cyanopropylphenyl



1. Toluene	110°C
2. Hexanol	158°C
3. Phenol	181°C
4. Decane (C10)	174°C
5. Naphthalene	218°C
6. Dodecane (C12)	216°C

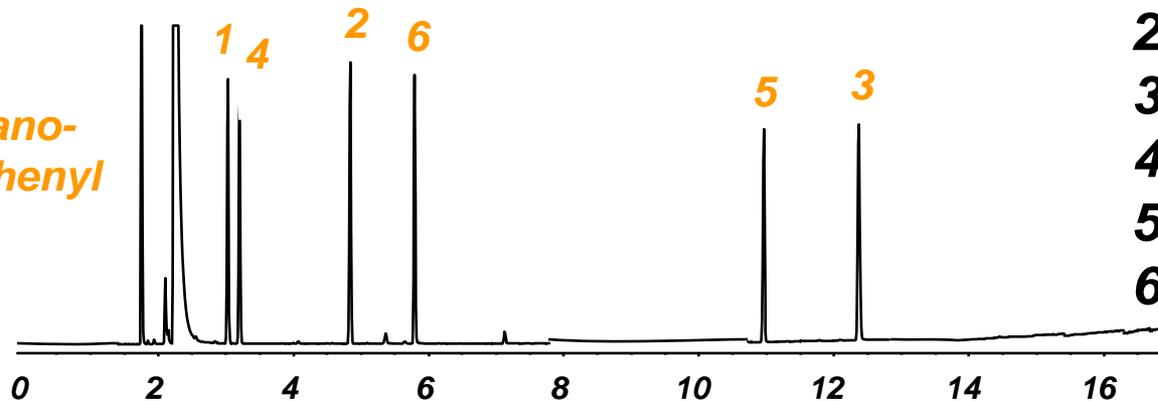


Strong Dispersion
None/Strong Dipole (Ph/CNPr)
Weak/Moderate H Bonding (Ph/CNPr)

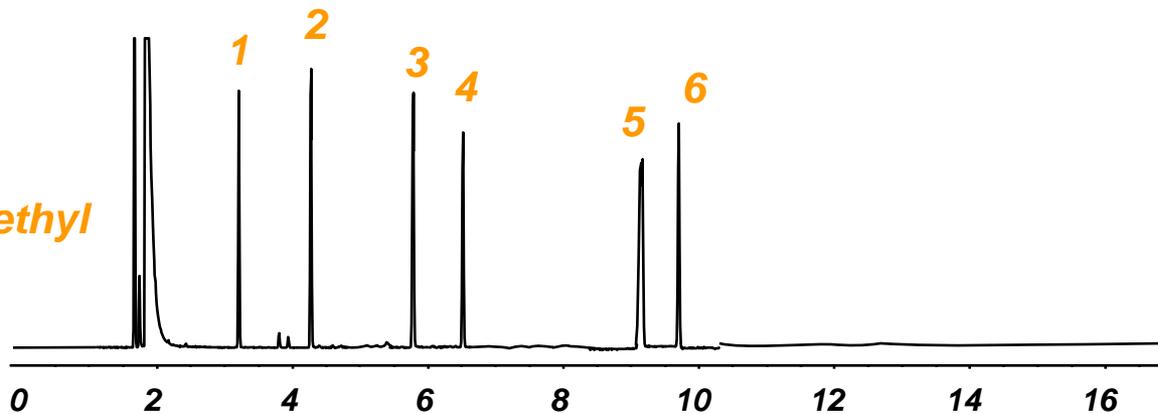
50% Cyanopropylphenyl

1. Toluene
2. Hexanol
3. Phenol
4. Decane (C10)
5. Naphthalene
6. Dodecane (C12)

50% Cyano-
propylphenyl

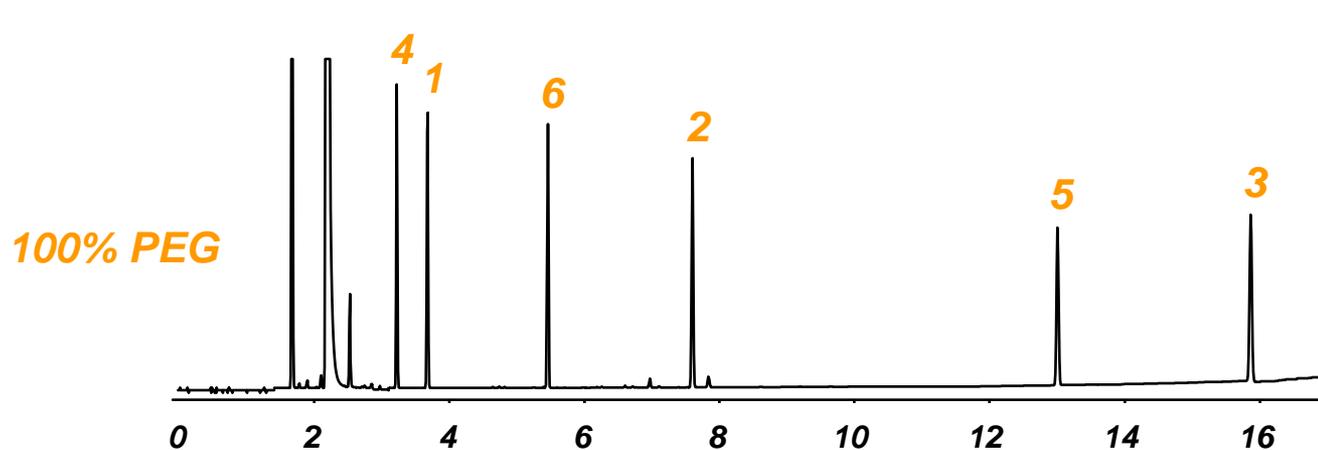


100% Methyl

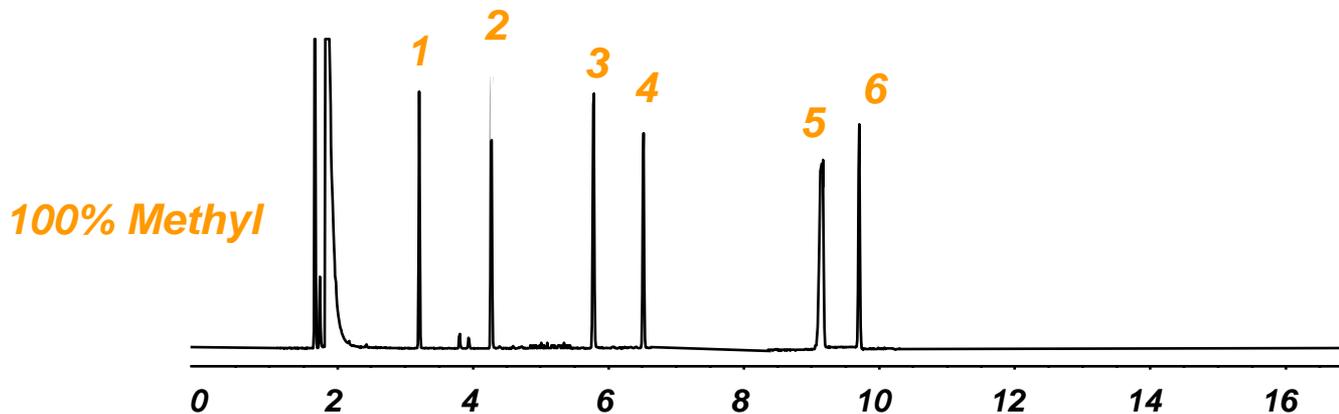


Strong Dispersion
None/Strong Dipole (Ph/CNPr)
Weak/Moderate H Bonding (Ph/CNPr)

100% Polyethylene Glycol



1. Toluene
2. Hexanol
3. Phenol
4. Decane (C10)
5. Naphthalene
6. Dodecane (C12)



Strong Dispersion
Strong Dipole
Moderate H Bonding

Stationary Phase Selection

Part 1

- Existing Information
- Selectivity
- Polarity
- Critical Separations
- Temperature Limits

Stationary Phase Selection

Part 2

- Capacity
- Analysis Time
- Bleed
- Versatility
- Selective Detectors

Column Dimensions

- Inner Diameter
- Length
- Film Thickness

Column Diameter

Capillary Columns

I.D. (mm)	Common Name
0.53	Megabore
0.45	High speed Megabore
0.32	Wide
0.20-0.25	Narrow
0.18	Minibore

Column Diameter

Theoretical Efficiency

I.D. (mm)	N/m
0.10	11905
0.18	6666
0.20	5941
0.25	4762
0.32	3717
0.53	2242

k = 5

Efficiency and Resolution Relationship

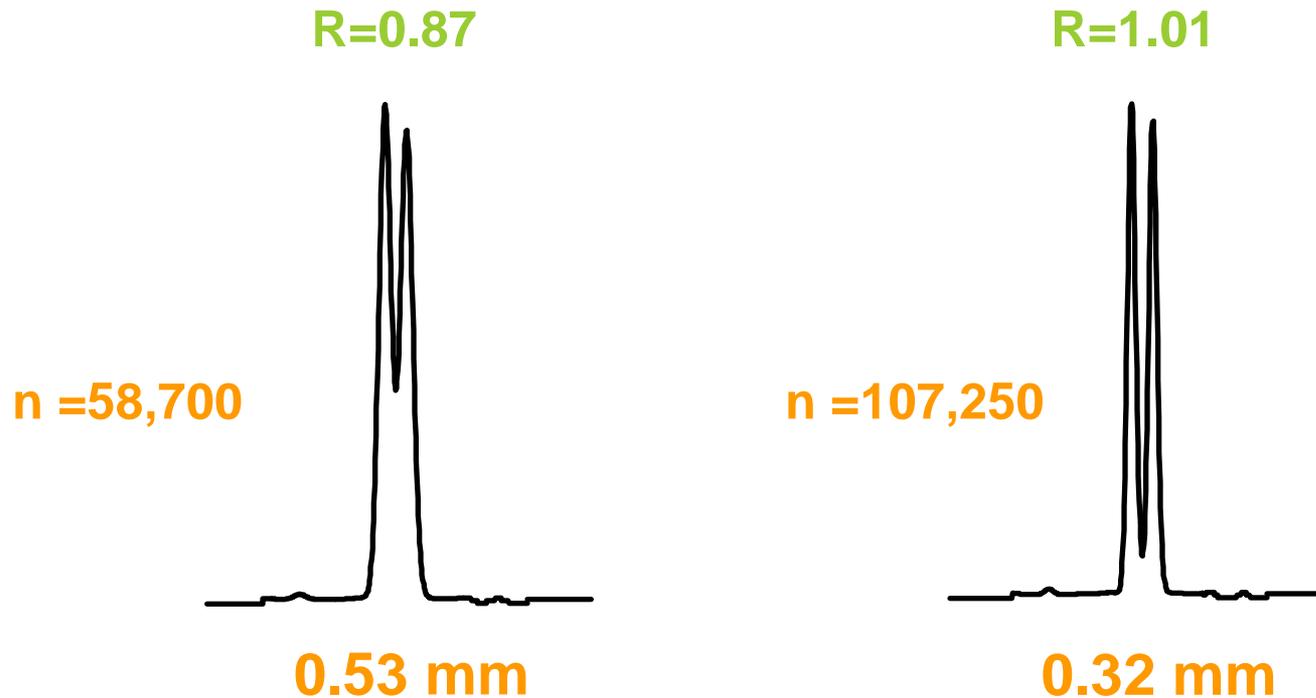
$$\sqrt{N} \propto R_s$$

Efficiency X 4 = Resolution X 2

Column Diameter

Resolution

180°C isothermal



Square root of resolution is inversely proportional to column diameter

Column Diameter

Inlet Head Pressures

Helium

I.D (mm)	Pressure (psig)
0.10	225-250
0.20	25-35
0.25	15-25
0.32	10-20
0.53	2-4

30 meters

Hydrogen pressures x 1/2

Column Diameter

Capacity

Like Polarity Phase/Solute

I.D. (mm)	Capacity (ng)
0.20	50-100
0.25	75-150
0.32	125-250
0.53	200-400

0.25 μm film thickness

Column Diameter

Carrier Gas Flow Rate

- Smaller diameters for low flow situations (e.g., GC/MS)
- Larger diameters for high flow situations (e.g., purge & trap, headspace, gas sample valve)

Column Length

Most common: 15-60 meters

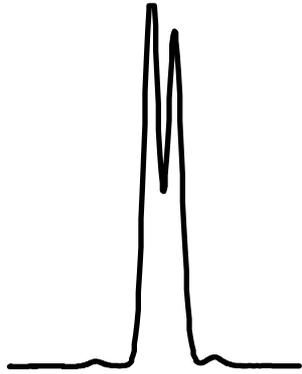
Available: 5-150 meters

Column Length

Resolution and Retention

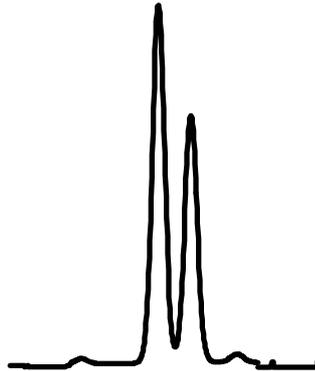
210°C isothermal

R=0.84
2.29 min



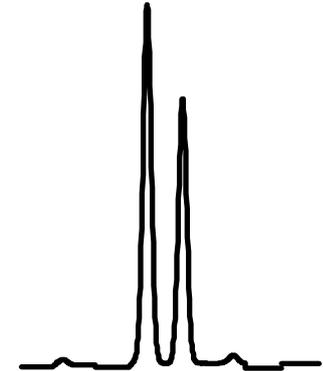
15 m

R=1.16
4.82 min



30 m

R=1.68
8.73 min

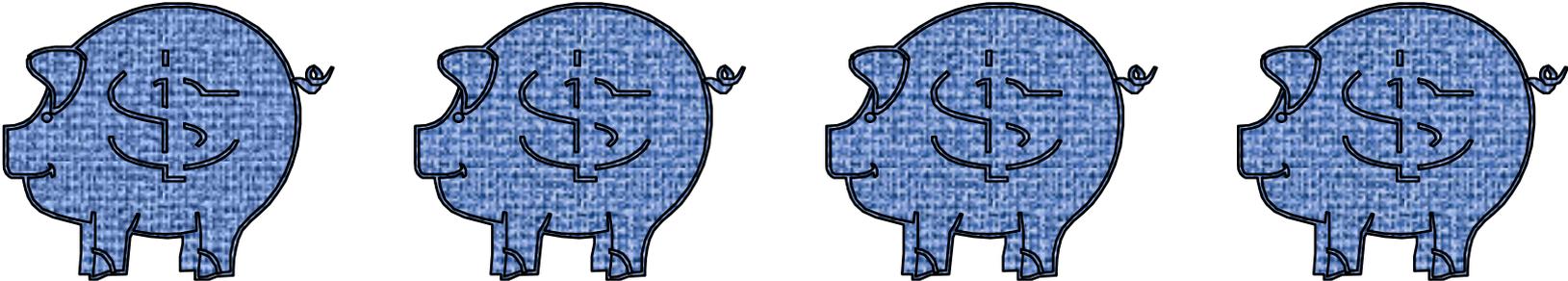
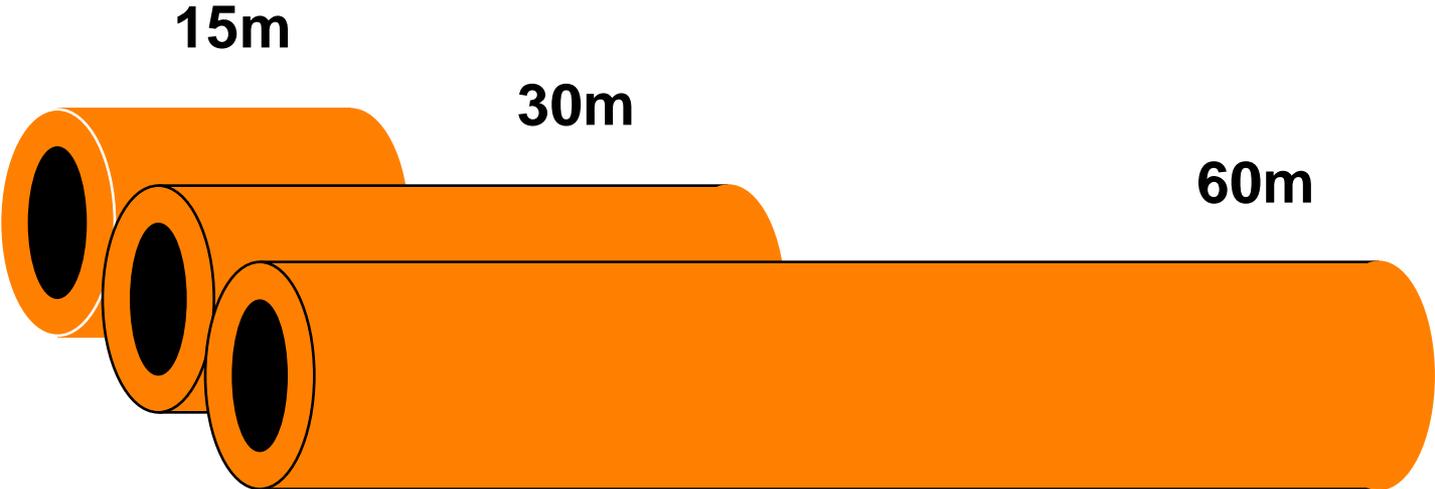


60 m

Resolution is proportional to the square root of column length
Isothermal: Retention is proportional to length
Temperature program: 1/3-1/2 of isothermal values

Column Length

Cost



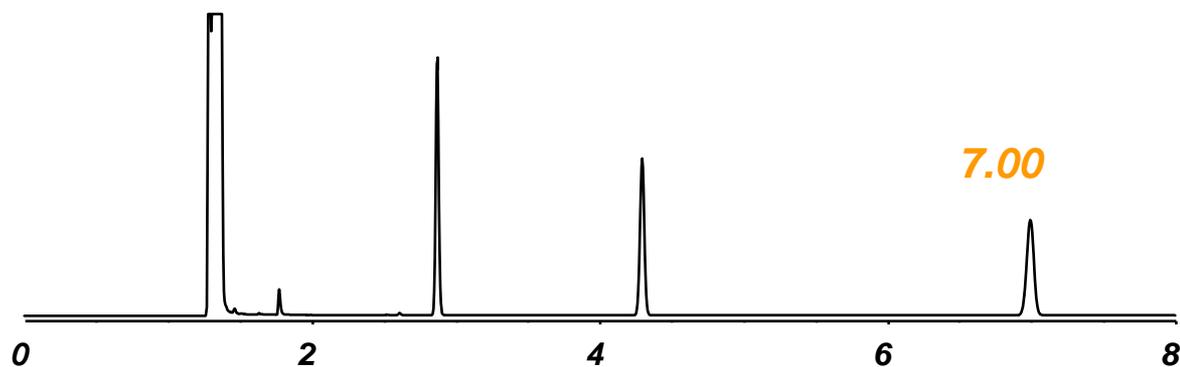
Film Thickness

- Most common: 0.1-3.0 μm
- Available: 0.1-10.0 μm

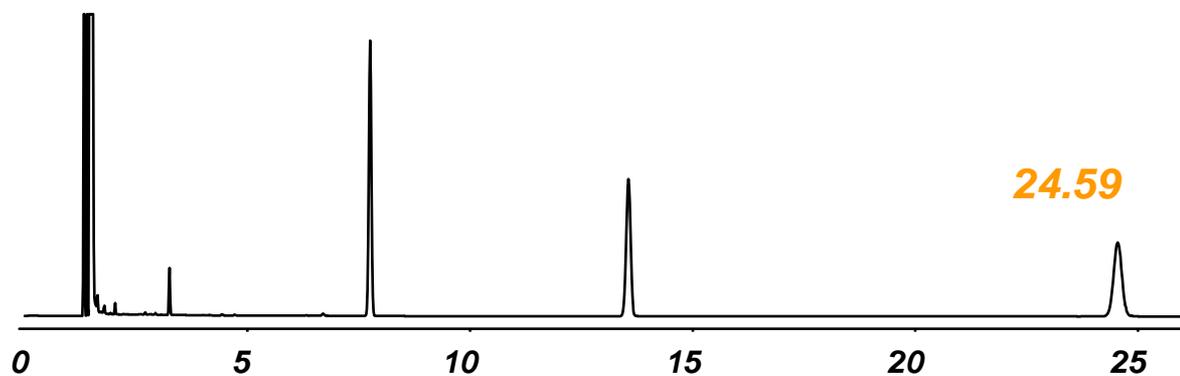
Film Thickness

Retention

100°C Isothermal



0.25 μm

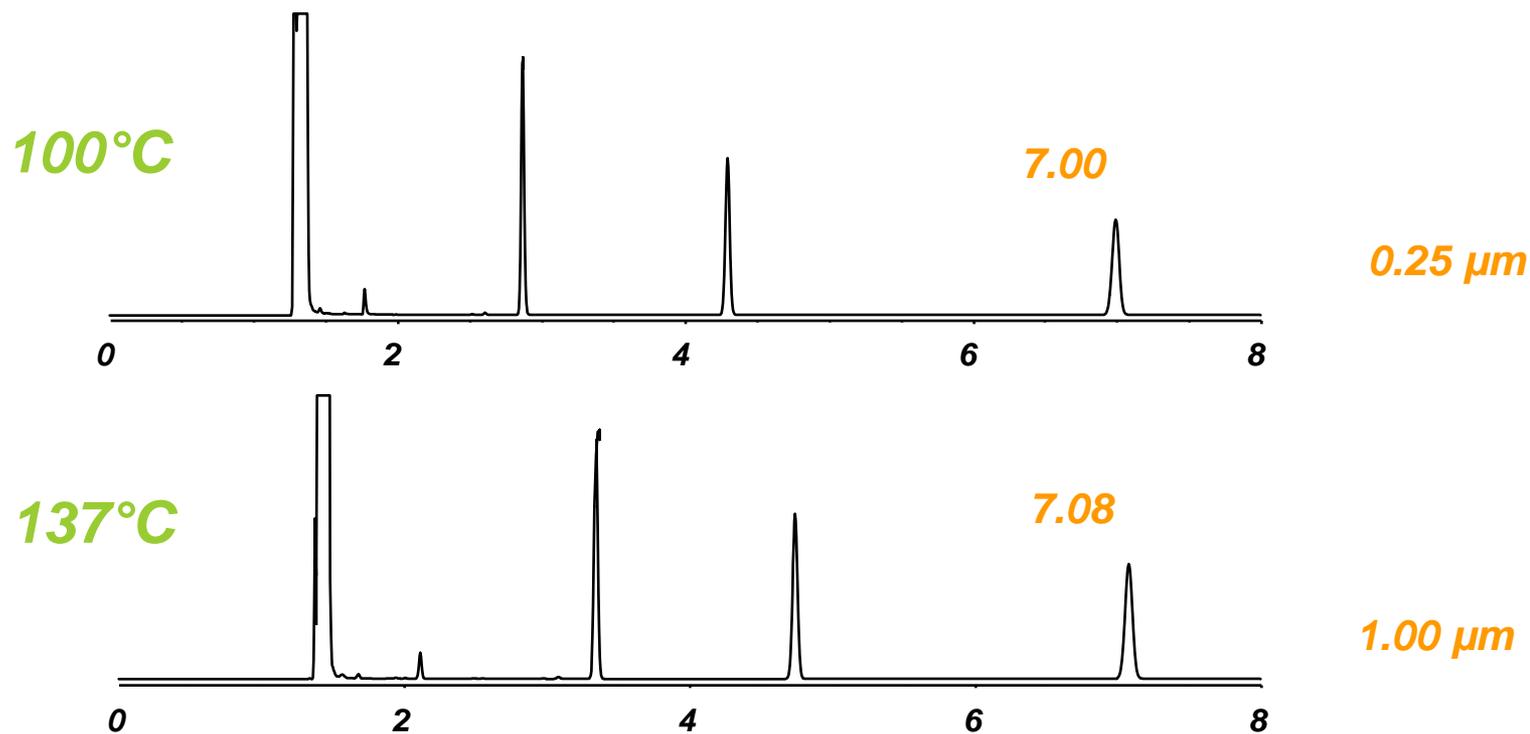


1.00 μm

Isothermal: Retention is proportional to film thickness
Temperature program: 1/3-1/2 of isothermal values

Film Thickness

Equal Retention: Isothermal

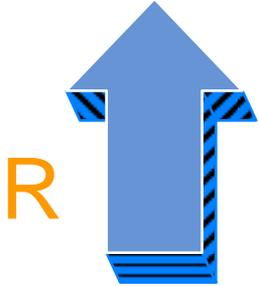


DB-1, 30 m x 0.32 mm ID
He at 37 cm/sec
C10, C11, C12

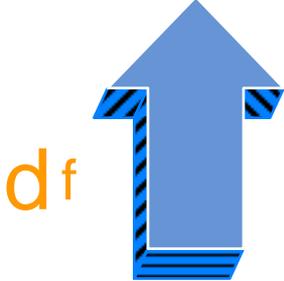
Film Thickness

Resolution

When solute $k < 5$

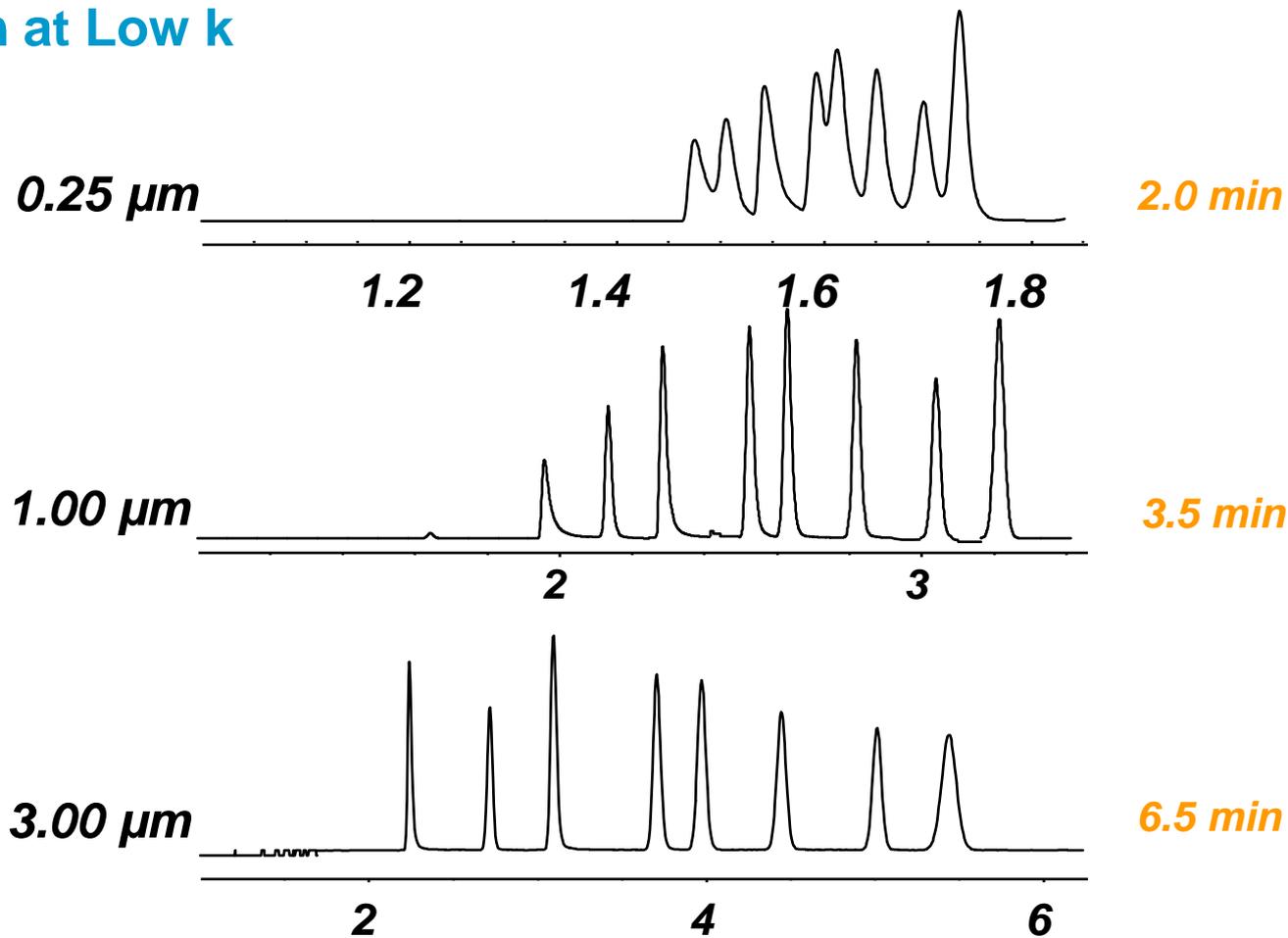


When solute $k > 5$



Film Thickness

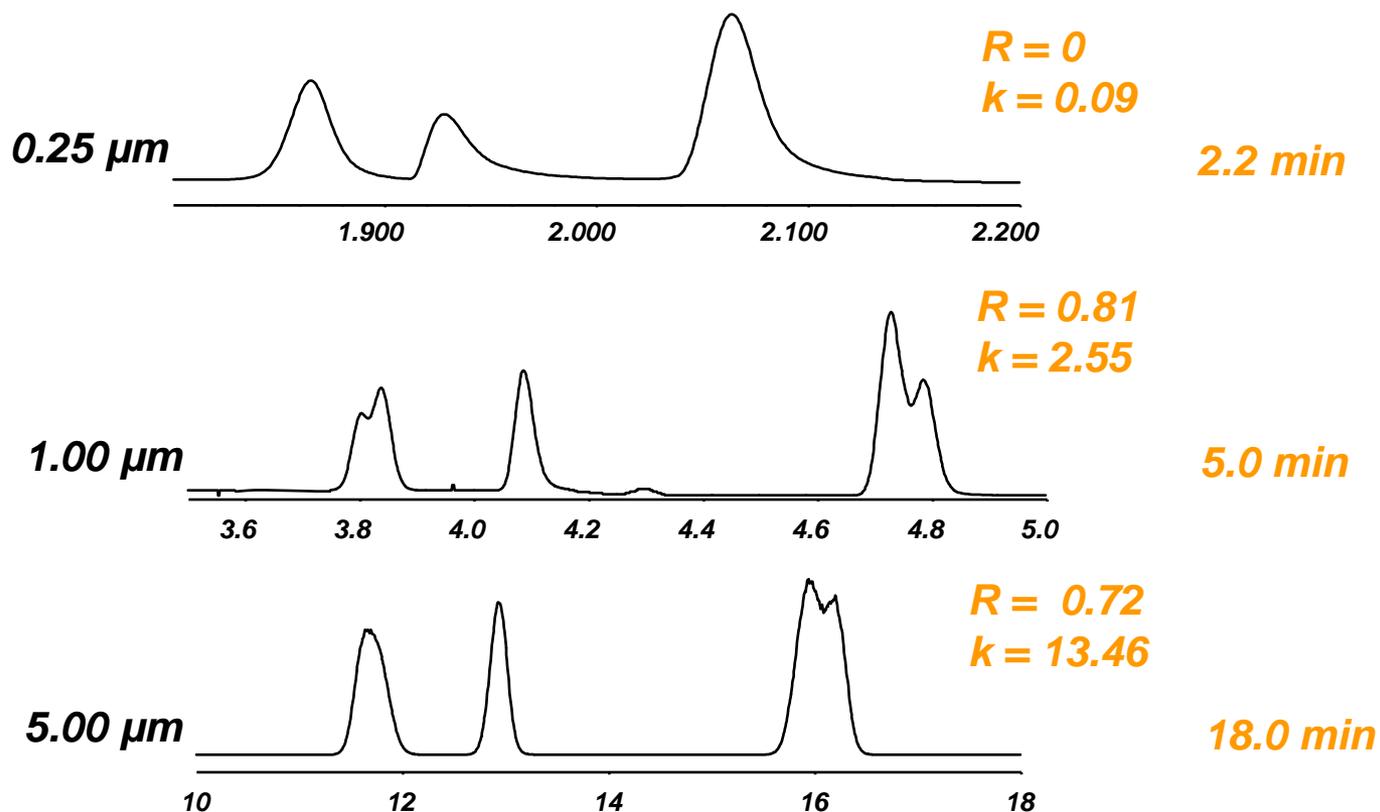
Resolution at Low k



DB-1, 30 m x 0.32 mm ID
40°C isothermal, He at 35 cm/sec
Solvent mixture

Film Thickness

Resolution at High k



DB-1, 30 m x 0.32 mm ID
40°C isothermal, He at 35 cm/sec
Solvent mixture

Film Thickness

Capacity

Like Polarity Phase/Solute

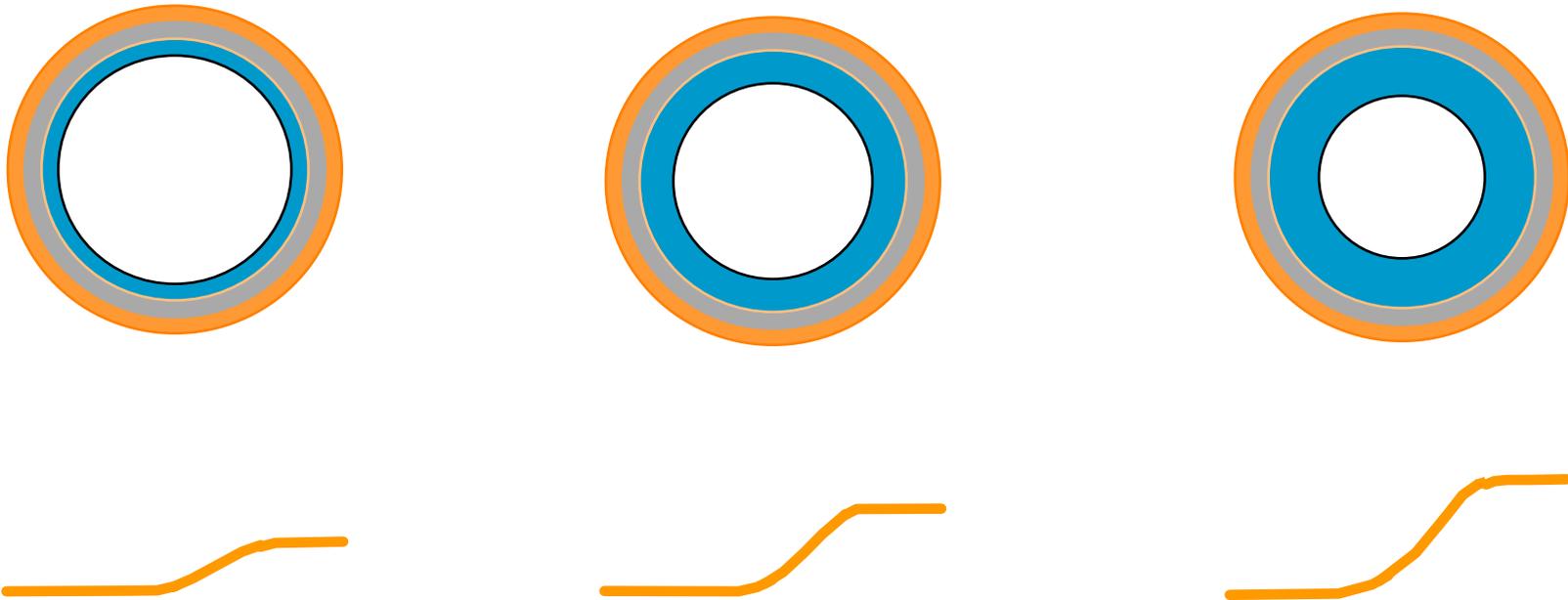
Thickness (um)	Capacity (ng)
0.10	50-100
0.25	125-250
1.0	500-1000
3.0	1500-3000
5.0	2500-5000

0.32 mm I.D.

Film Thickness

Bleed

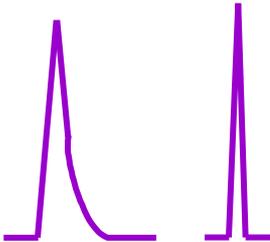
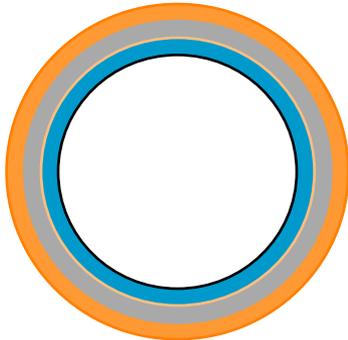
More stationary phase = More degradation products



Film Thickness

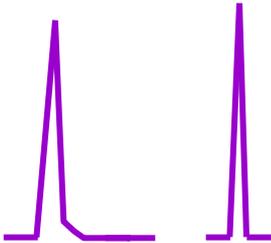
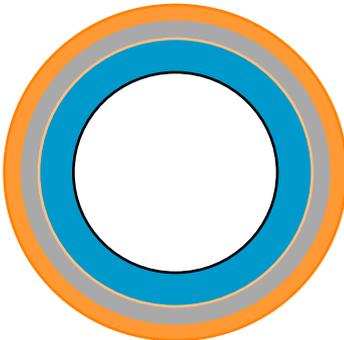
Inertness Summary

0.25



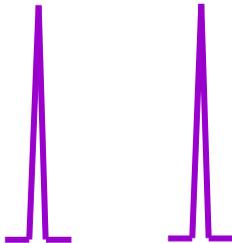
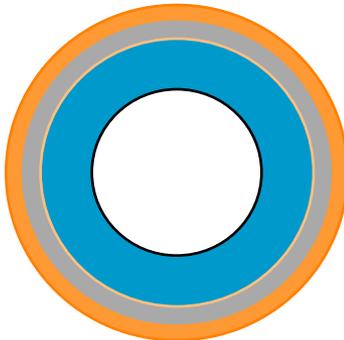
active inactive

1.0



active inactive

3.0



active inactive

Column Dimensions

Diameter Summary

To Increase	Make Diameter
Resolution	Smaller
Retention	Smaller
Pressure	Smaller
Flow rate	Larger
Capacity	Larger

Column Dimensions

Length Summary

To Increase

Make Length

Resolution

Longer

Retention

Longer

Pressure

Longer

Cost

Longer

Column Dimensions

Film Thickness Summary

To Increase

Make Film

Retention

Thicker

Resolution ($k < 5$)

Thicker

Resolution ($k > 5$)

Thinner

Capacity

Thicker

Inertness

Thicker

Bleed

Thicker

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