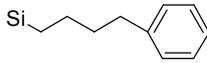
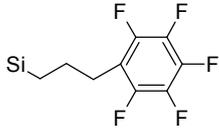


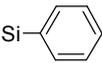
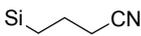
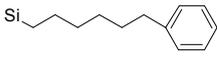
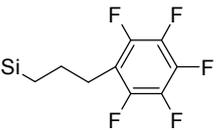
Comparison of separation selectivity among YMC-Triart series

**Separation of structural isomers by utilizing unique
separation characteristics of Triart Phenyl/PFP**

Specifications of Triart series for reversed phase chromatography

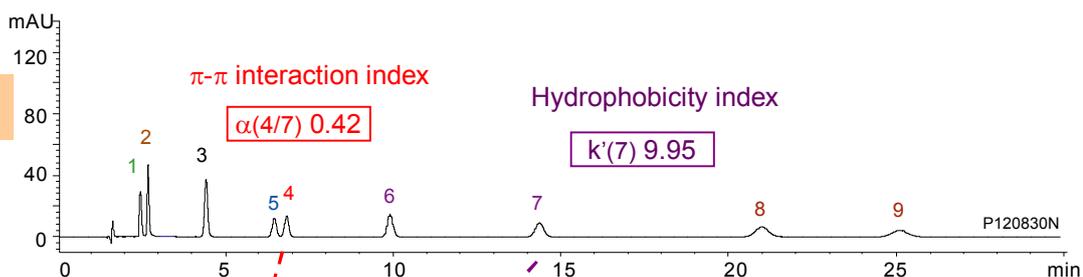
	Triart C18	Triart C8	Triart Phenyl	Triart PFP
Functional group	$\text{Si}-\text{C}_{18}\text{H}_{37}$	$\text{Si}-\text{C}_8\text{H}_{17}$		
Base	organic/inorganic hybrid silica			
Particle	5 μm , 3 μm , 1.9 μm			
Pore	12 nm			
Endcapping	Multi-stage endcapping			No
pH range	1~12		1~10	1~8
Upper limit of temperature	70°C for pH 1~7, 50°C for pH 7~12		50°C	

Specifications of other columns

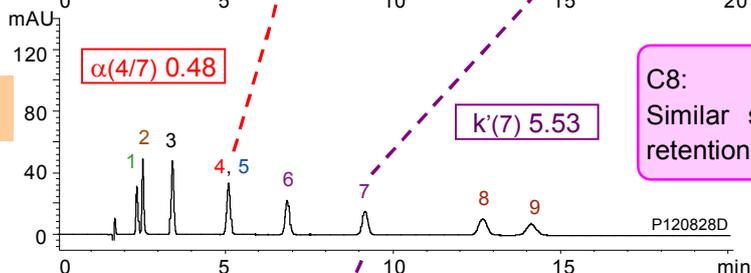
	YMC-Pack Ph	YMC-Pack CN	Conventional Phenyl-Hexyl	Conventional PFP
Functional group				
Base	Silica	Silica	Silica	Silica
Particle	5 μm , 3 μm	5 μm , 3 μm	5 μm , 3.5 μm , 1.8 μm	5 μm , 3 μm
Pore	12 nm	12 nm	9.5 nm	12 nm
Endcapping	Yes	Yes	Yes	Yes
pH range	2~7.5	2~7.5	2~8	2~7.5
Upper limit of temperature	50°C	50°C	60°C	75°C

Comparison of Separation Selectivity Among YMC Reversed Phase Columns

Triart C18

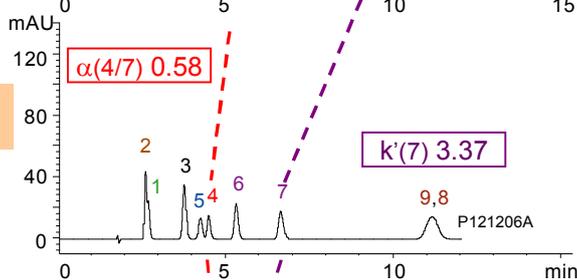


Triart C8



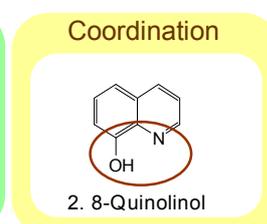
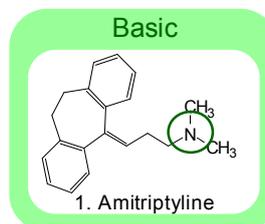
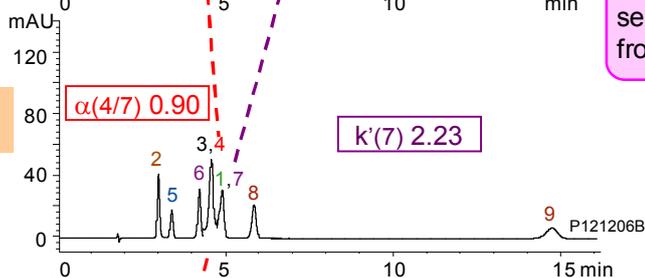
C8:
Similar selectivity to C18 and shorter retention times than C18

Triart Phenyl

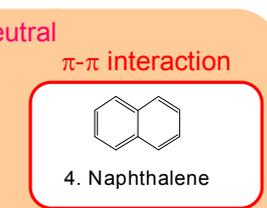
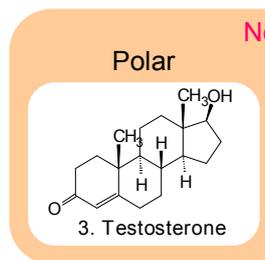
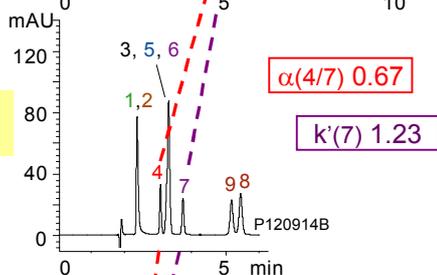


Phenyl and PFP:
 π - π interaction and polar interaction as well as hydrophobic interaction contribute to separation. Different separation selectivity from Triart C18/C8.

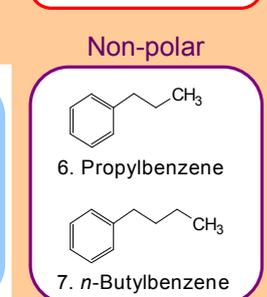
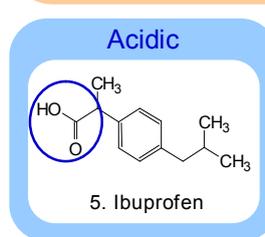
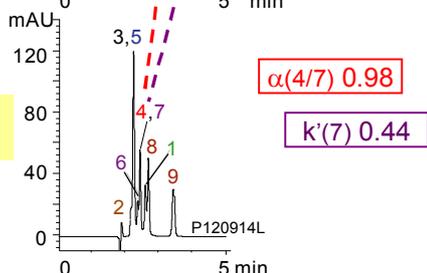
Triart PFP



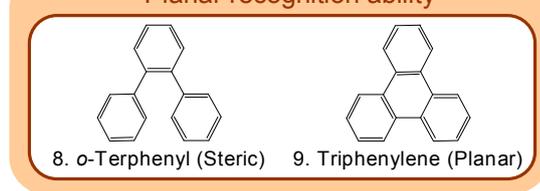
YMC-Pack Ph



YMC-Pack CN

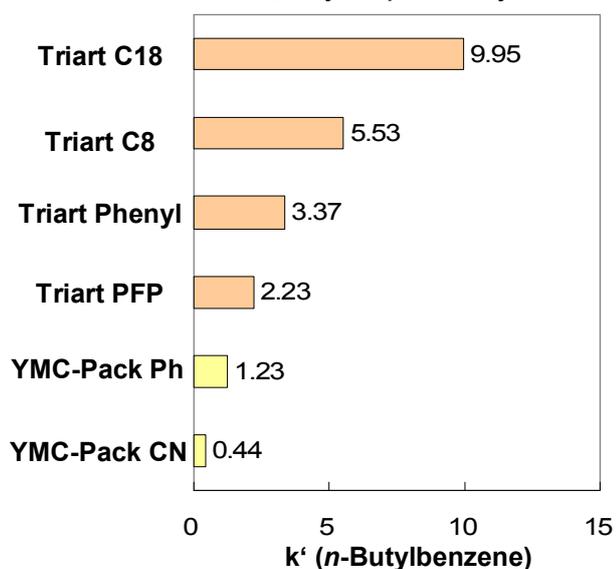


Planar recognition ability

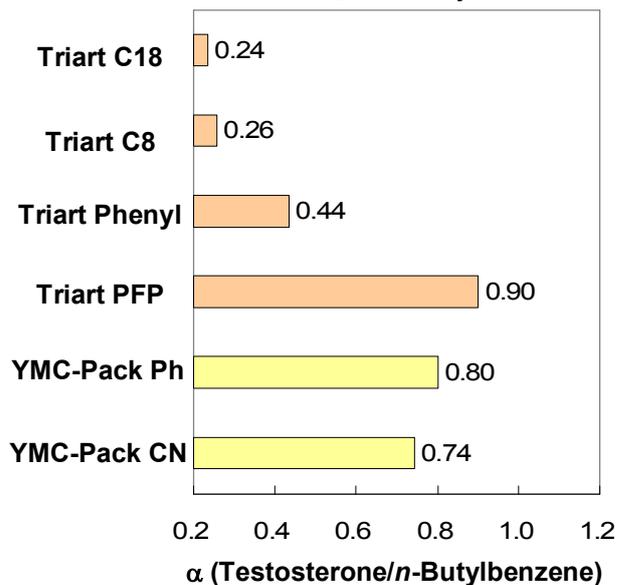


Column : 5 μ m, 150 X 3.0 mm I.D.
 Eluent : 20 mM KH_2PO_4 - H_3PO_4 (pH3.1)/methanol (25/75)
 Flow rate : 0.425 mL/min
 Temperature : 40°C
 Detection : UV at 265 nm
 Injection : 4 μ L

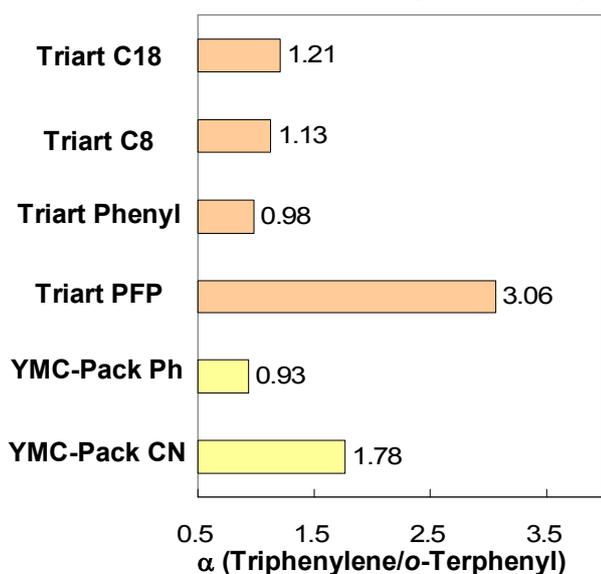
① Hydrophobicity



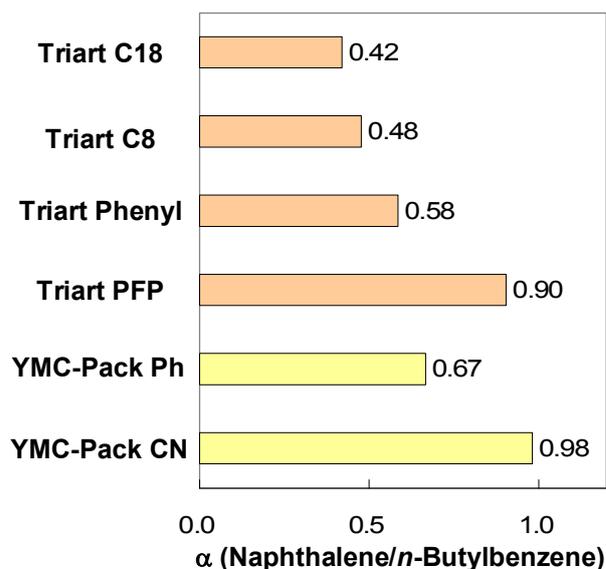
② Polarity



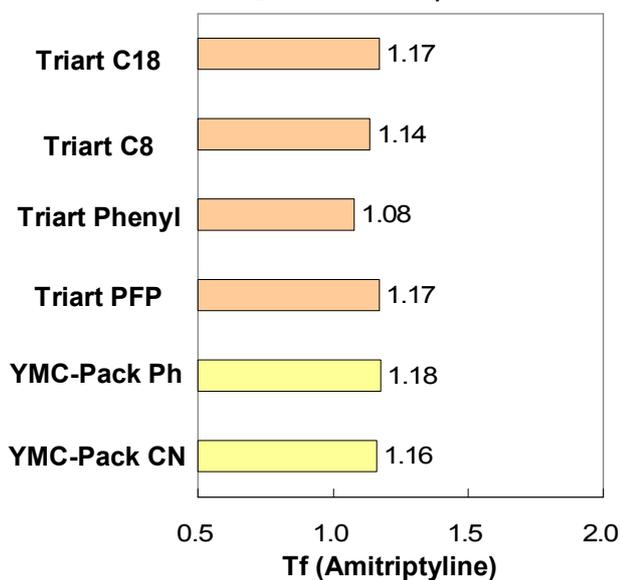
③ Planar recognition ability



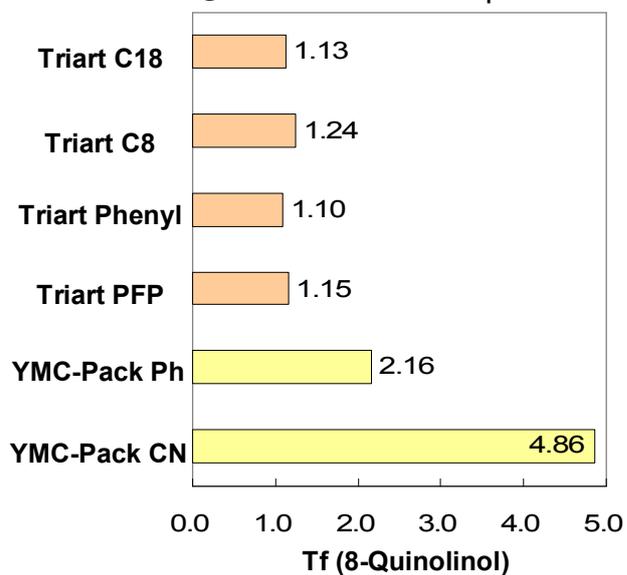
④ π - π interaction



⑤ Basic compound



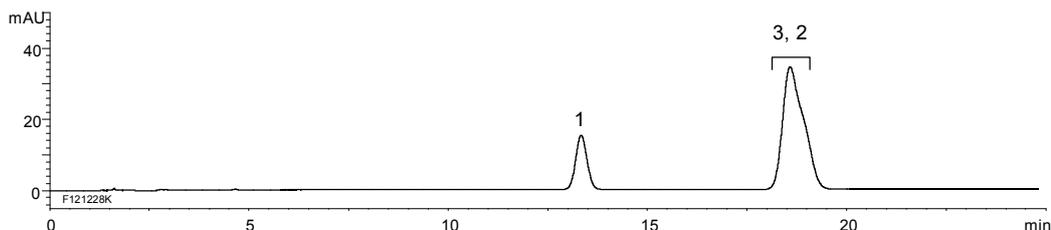
⑥ Coordination compound



Separation Comparison of Terphenyl Positional Isomers

Triart C18

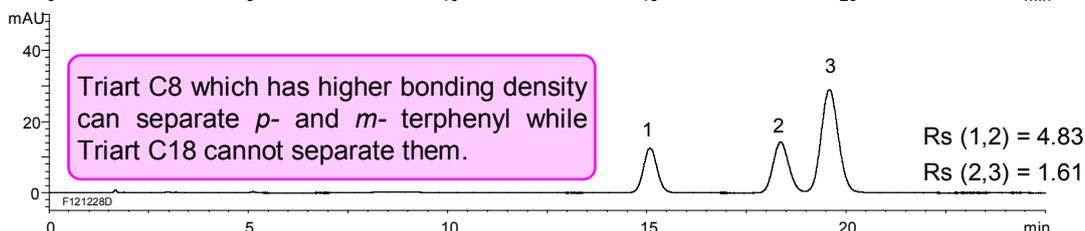
80% methanol



Triart C8

75% methanol

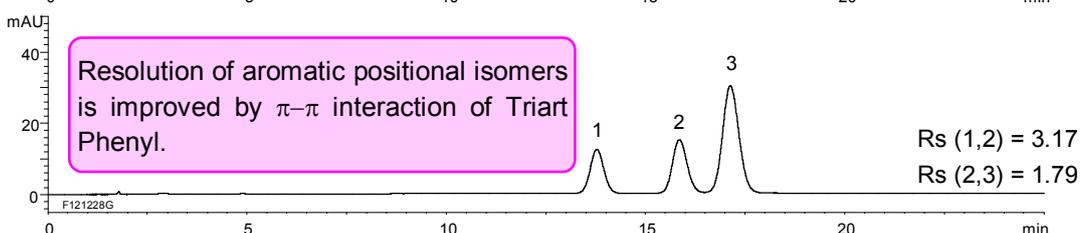
Triart C8 which has higher bonding density can separate *p*- and *m*- terphenyl while Triart C18 cannot separate them.



Triart Phenyl

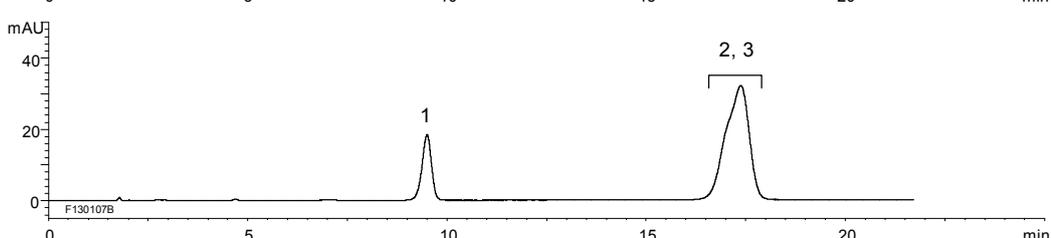
75% methanol

Resolution of aromatic positional isomers is improved by π - π interaction of Triart Phenyl.



Triart PFP

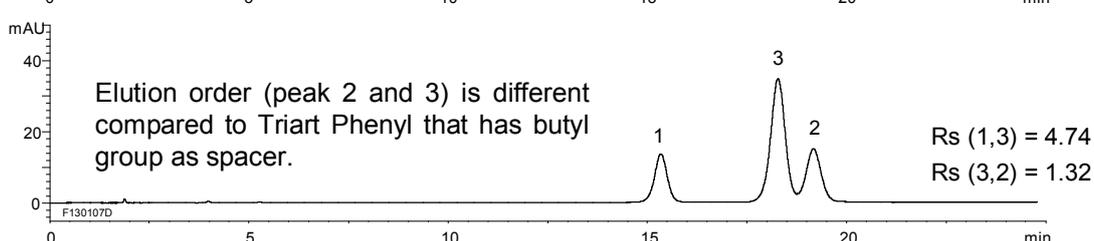
70% methanol



YMC-Pack Ph

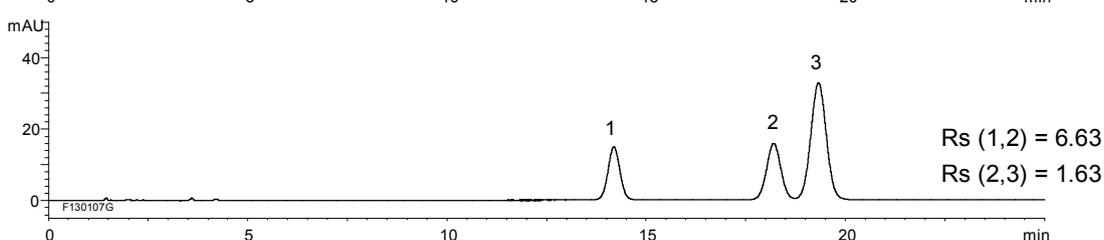
65% methanol

Elution order (peak 2 and 3) is different compared to Triart Phenyl that has butyl group as spacer.



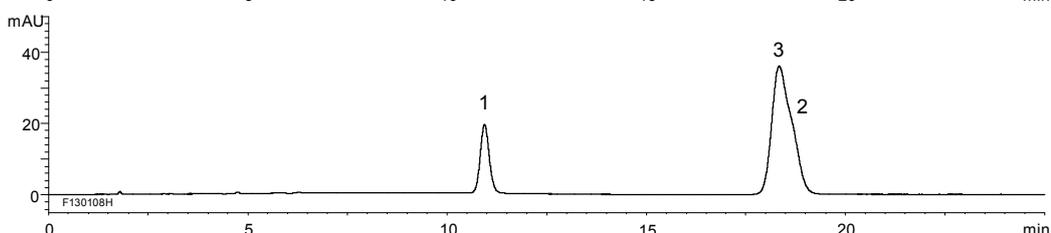
Conventional Phenyl-Hexyl

75% methanol



Conventional PFP

70% methanol



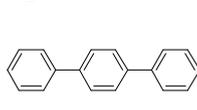
Terphenyl positional isomers

1



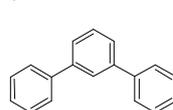
o-Terphenyl

2



p-Terphenyl

3

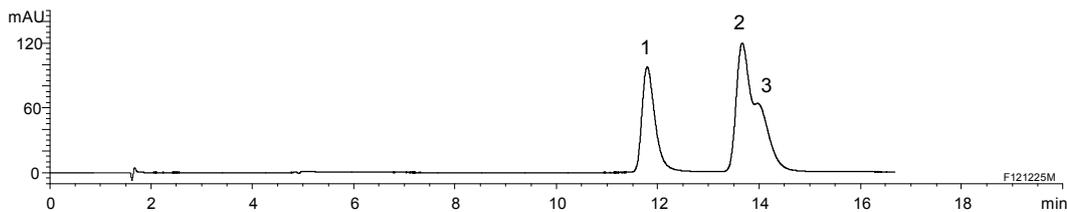


m-Terphenyl

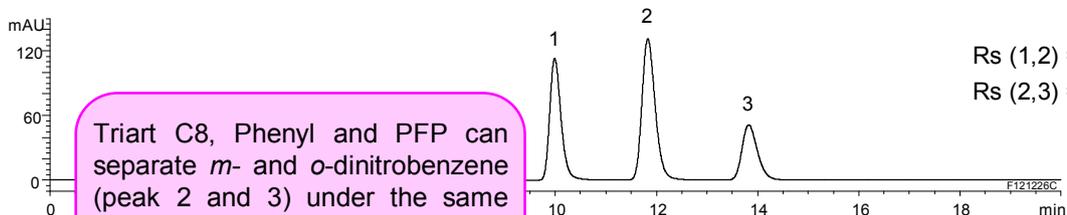
Column	: 5 μ m, 150 X 3.0 or 4.6 mmI.D.
Eluent	: methanol/water
Flow rate	: 0.425 mL/min for 3.0 mmI.D. 1.0 mL/min for 4.6 mmI.D.
Temperature	: 30°C
Detection	: UV at 254 nm
Injection	: 2.0 μ L for 3.0 mmI.D. 4.7 μ L for 4.6 mmI.D.

Separation Comparison of Dinitrobenzene Positional Isomers

Triart C18



Triart C8

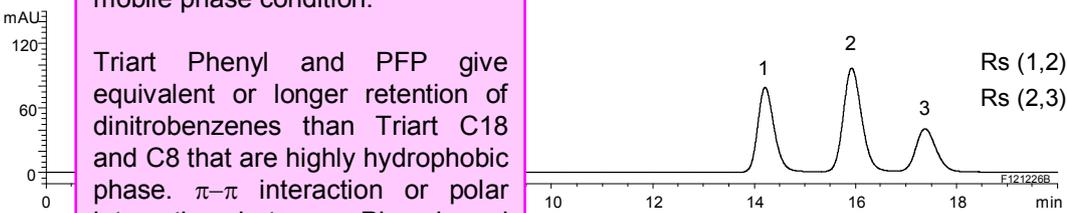


$R_s(1,2) = 4.29$

$R_s(2,3) = 3.93$

Triart C8, Phenyl and PFP can separate *m*- and *o*-dinitrobenzene (peak 2 and 3) under the same mobile phase condition.

Triart Phenyl

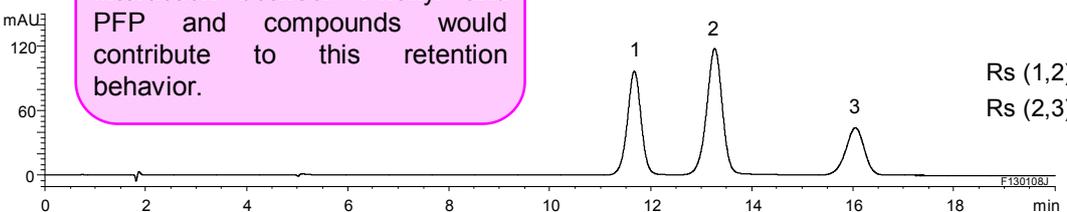


$R_s(1,2) = 2.79$

$R_s(2,3) = 2.11$

Triart Phenyl and PFP give equivalent or longer retention of dinitrobenzenes than Triart C18 and C8 that are highly hydrophobic phase. π - π interaction or polar interaction between Phenyl and PFP and compounds would contribute to this retention behavior.

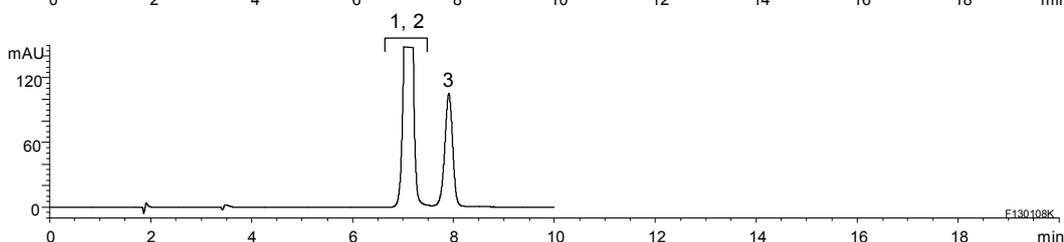
Triart PFP



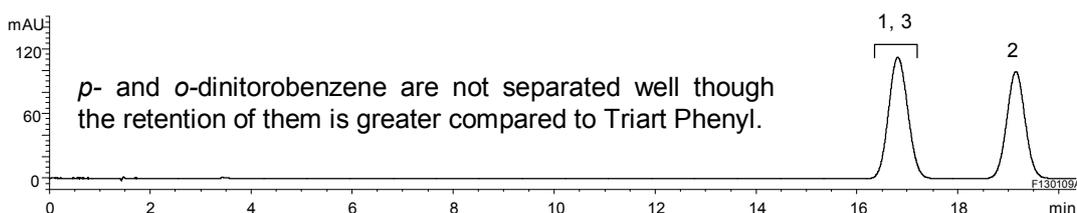
$R_s(1,2) = 3.11$

$R_s(2,3) = 4.62$

YMC-Pack Ph

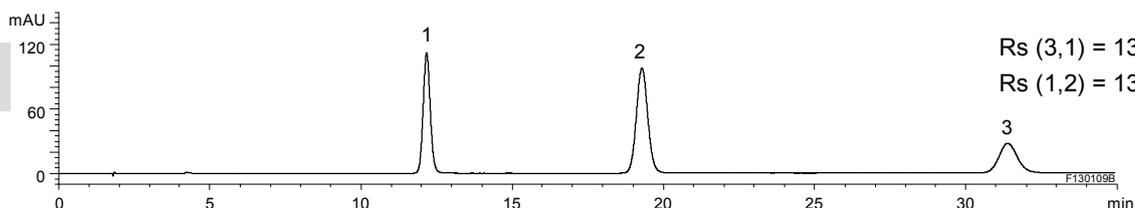


Conventional Phenyl-Hexyl



p- and *o*-dinitrobenzene are not separated well though the retention of them is greater compared to Triart Phenyl.

Conventional PFP

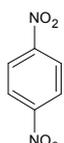


$R_s(3,1) = 13.09$

$R_s(1,2) = 13.80$

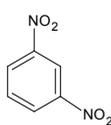
Dinitrobenzene positional isomers

1



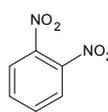
p-Dinitrobenzene

2



m-Dinitrobenzene

3

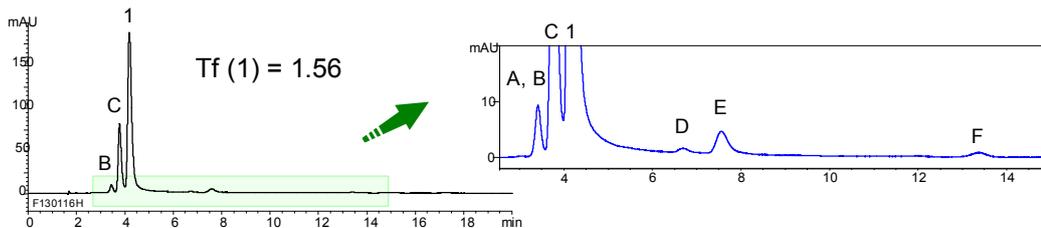


o-Dinitrobenzene

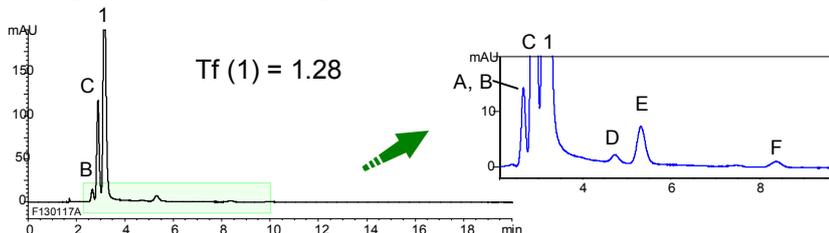
Column	: 5 μ m, 150 X 3.0 or 4.6 mm I.D.
Eluent	: methanol/water (40/60)
Flow rate	: 0.425 mL/min for 3.0 mm I.D. 1.0 mL/min for 4.6 mm I.D.
Temperature	: 30°C
Detection	: UV at 254 nm
Injection	: 2.1 μ L for 3.0 mm I.D. 5.0 μ L for 4.6 mm I.D.

Separation Comparison of Brilliant Blue FCF and Its Impurities

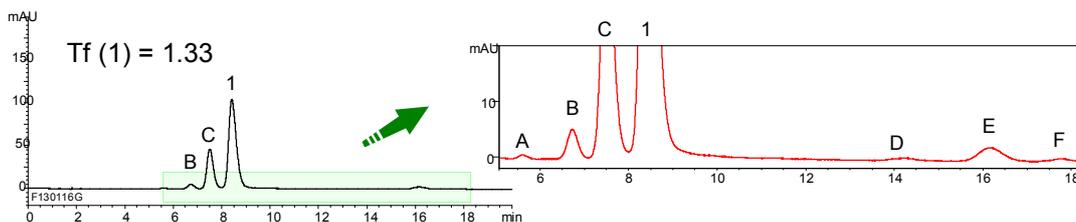
Triart C18
45% methanol



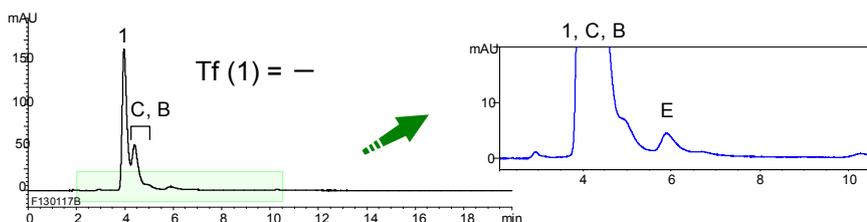
Triart C8
45% methanol



Triart Phenyl
45% methanol

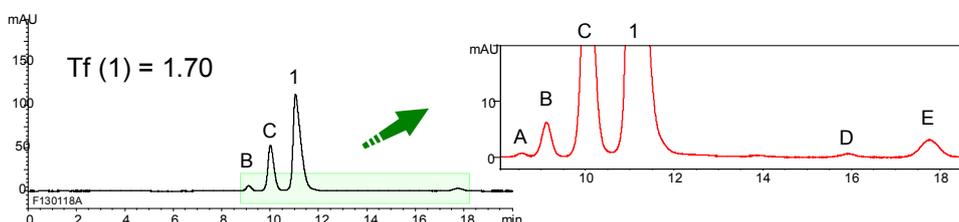


Triart PFP
45% methanol

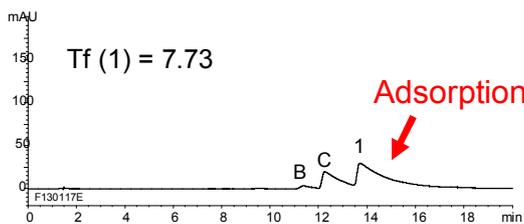


Elution order is different from that on phenyl or alkyl type stationary phase.

YMC-Pack Ph
55% methanol

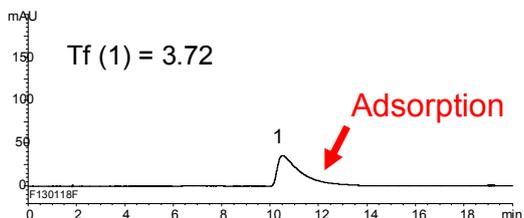


Conventional Phenyl-Hexyl
45% methanol

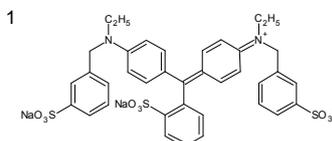


Phenyl type phases provide preferable retention and separation on brilliant blue FCF, acidic triphenylmethane dye, and its impurities.

Conventional PFP
45% methanol



Brilliant blue FCF



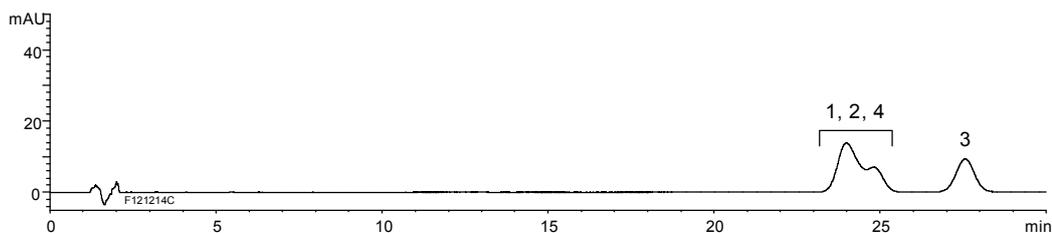
A - E
: Impurities in Brilliant blue FCF reagent

Brilliant blue FCF

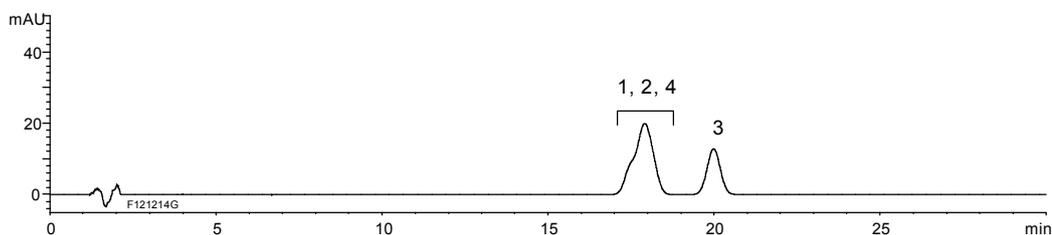
Column	: 5 μ m, 150 X 3.0 or 4.6 mmI.D.
Eluent	: 0.1% phosphoric acid/methanol
Flow rate	: 0.425 mL/min for 3.0 mmI.D. 1.0 mL/min for 4.6 mmI.D.
Temperature	: 40°C
Detection	: UV at 630 nm
Injection	: 2.0 μ L for 3.0 mmI.D. 4.7 μ L for 4.6 mmI.D.

Separation Comparison of Piperine *cis-trans* Isomers

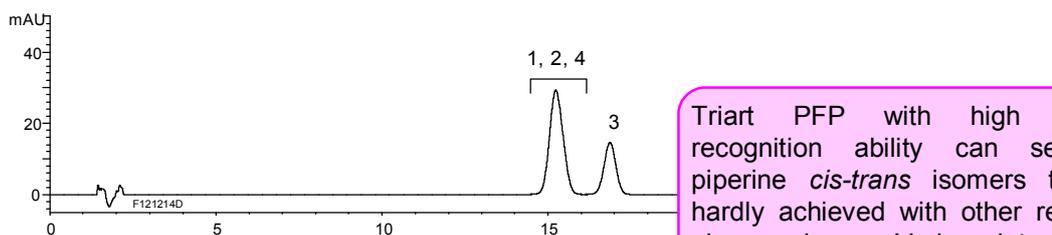
Triart C18



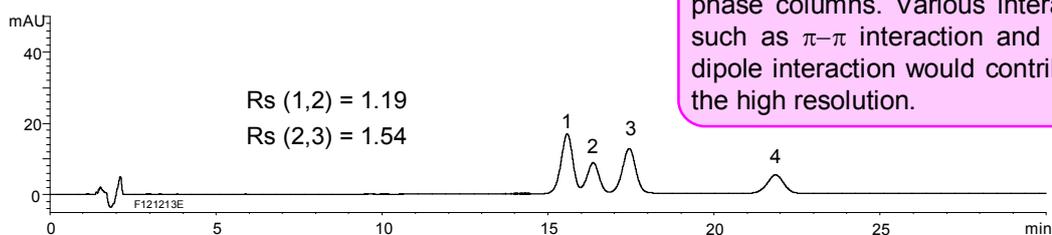
Triart C8



Triart Phenyl

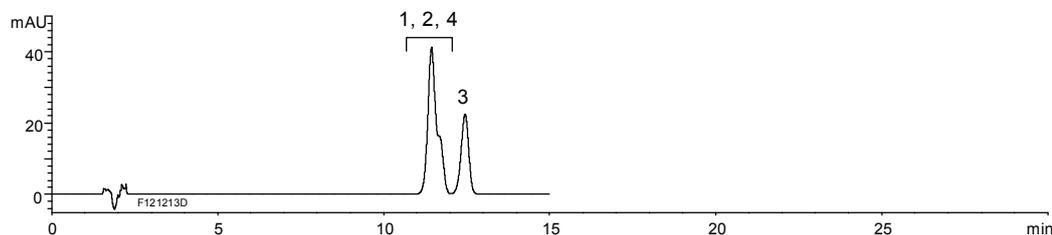


Triart PFP

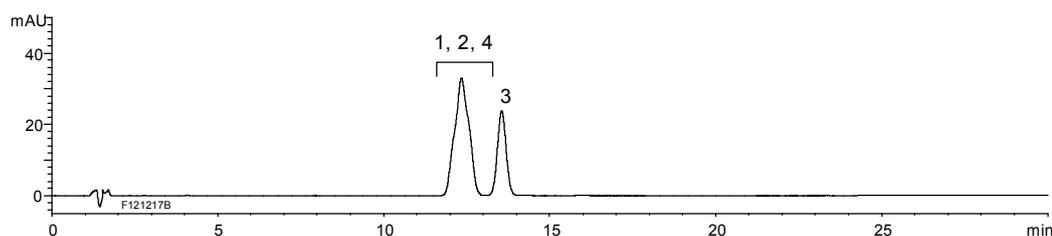


Triart PFP with high planar recognition ability can separate piperine *cis-trans* isomers that is hardly achieved with other reverse-phase columns. Various interactions such as $\pi-\pi$ interaction and dipole-dipole interaction would contribute to the high resolution.

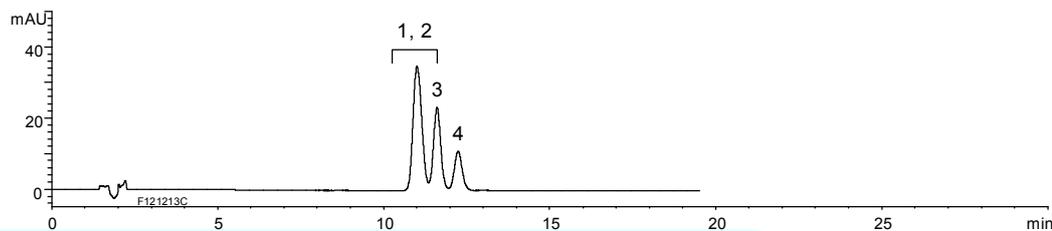
YMC-Pack Ph



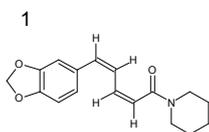
Conventional Phenyl-Hexyl



Conventional PFP

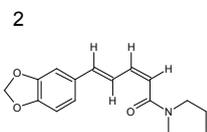


Piperine and its isomers



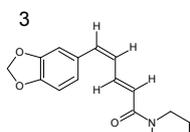
Chavicine

Z,Z-(*cis-cis*)-piperine



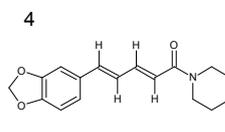
Isopiperine

Z,E-(*cis-trans*)-piperine



Isochavicine

E,Z-(*trans-cis*)-piperine



Piperine

E,E-(*trans-trans*)-piperine

Column	: 5 μ m, 150 X 3.0 or 4.6 mm I.D.
Eluent	: 0.1% formic acid/acetonitrile (60/40)
Flow rate	: 0.425 mL/min for 3.0 mm I.D. 1.0 mL/min for 4.6 mm I.D.
Temperature	: 25°C
Detection	: UV at 280 nm
Injection	: 4.25 μ L for 3.0 mm I.D. 10.0 μ L for 4.6 mm I.D.