



Photos: CERN

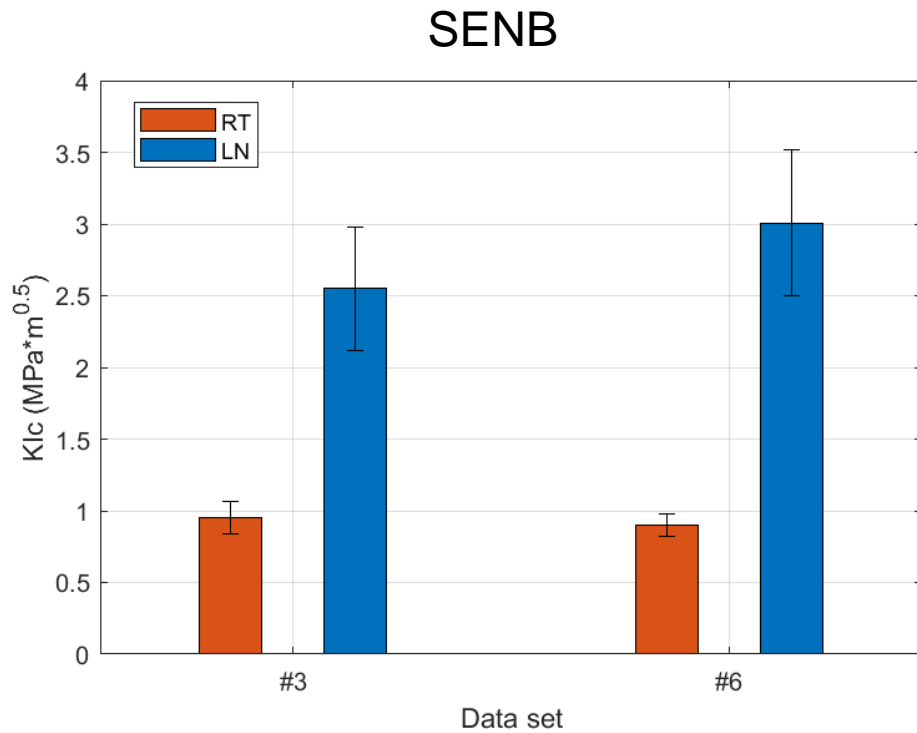
Tough epoxy systems

CHART Update
23.09.2021

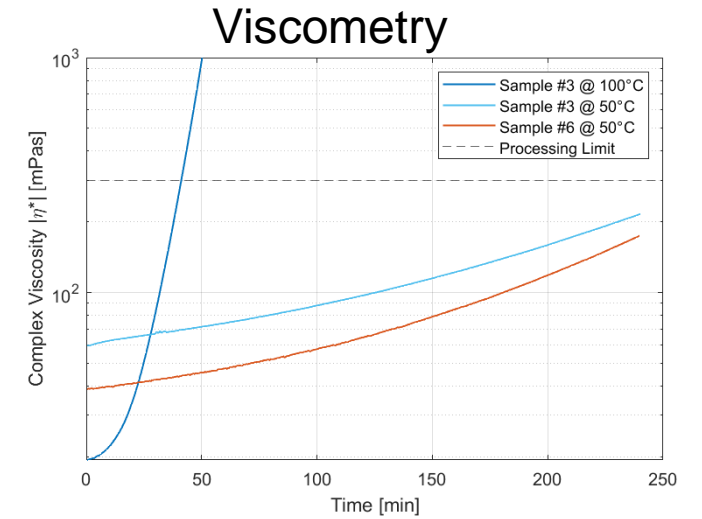
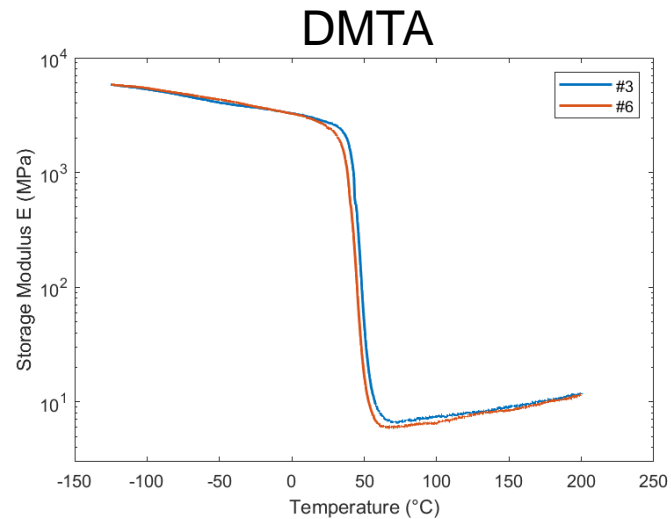
Pascal Studer
Laboratory for Soft Materials

Recap

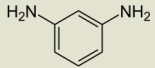
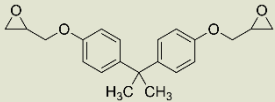
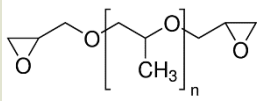
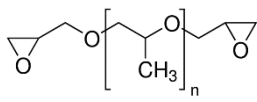
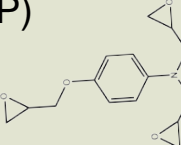
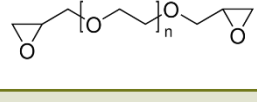
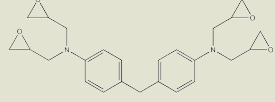
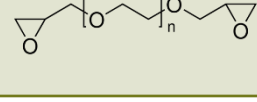

- Good toughness → Approaching MY750 values
- Pot life > 4h
- Low viscosity



#3 Wt. %	#6 Wt. %	Component
-	12.5	<chem>Nc1ccc(N)cc1</chem>
21	-	<chem>Nc1ccc(cc1)Cc2ccc(N)cc2</chem>
15.7	18.2	<chem>CC(C)(c1ccc(OCC2OC2)cc1)c3ccc(OCC4OC4)cc3</chem>
63.3	69.4	<chem>CC(C)OCC5OC5</chem>



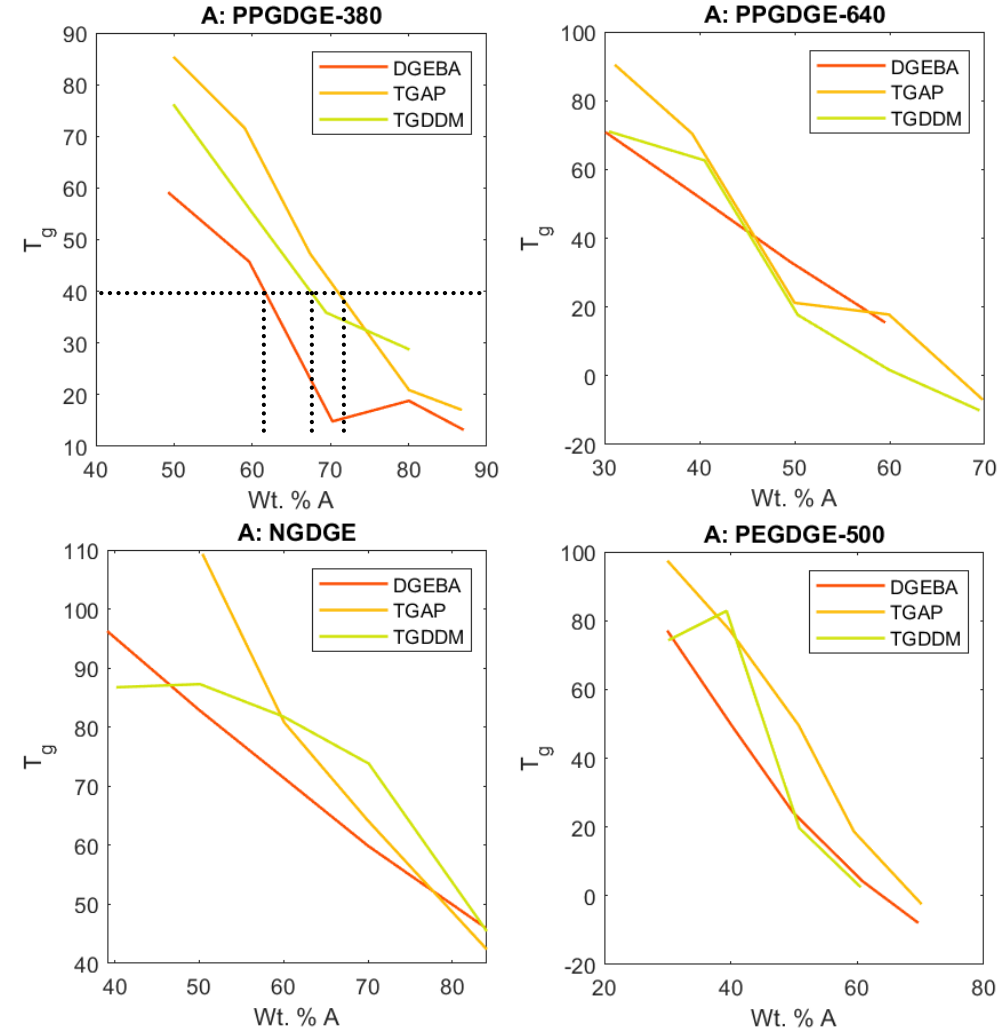
Systematic study – System & Objectives

Comp. C Amine	Comp. B Epoxide	Comp. A (f = 2) Epoxide, flexible chain
Meta-phenylene diamine (MPD) $f = 4$ 	Bisphenol A diglycidyl ether (DGEBA) $f = 2$ 	Poly(propylene glycol) diglycidyl ether 380 g/mol (PPGDGE-380) 
		Poly(propylene glycol) diglycidyl ether 640 g/mol (PPGDGE-640) 
	Triglycidyl p-aminophenol (TGAP) $f = 3$ 	Poly(ethylene glycol) diglycidyl ether 400 g/mol (PEGDGE-400) 
	Tetraglycidyl methylenedianiline (TGDDM) $f = 4$ 	Poly(ethylene glycol) diglycidyl ether 500 g/mol (PEGDGE-500) 
	Neopentyl glycol diglycidyl ether (NGDGE) 	

- Relate structural parameters (crosslink density, deformability,...) to toughness
- Each composition is a combination of A, B, C
 - Stoichiometric
 - One degree of freedom (I use **wt. % A**)
- Increase toughness even more
 - Maximize wt. % A, so that $T_g = 40\text{ °C}$ & see which is most effective

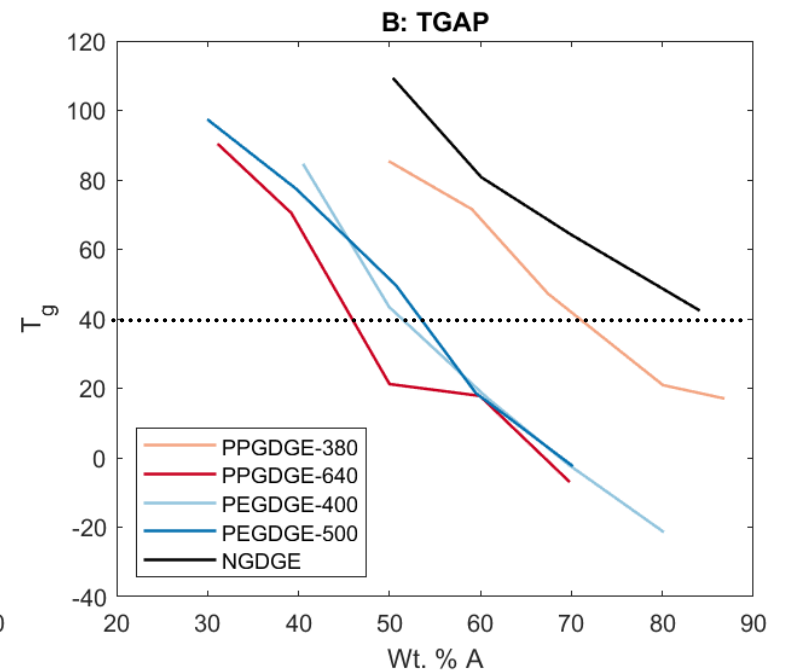
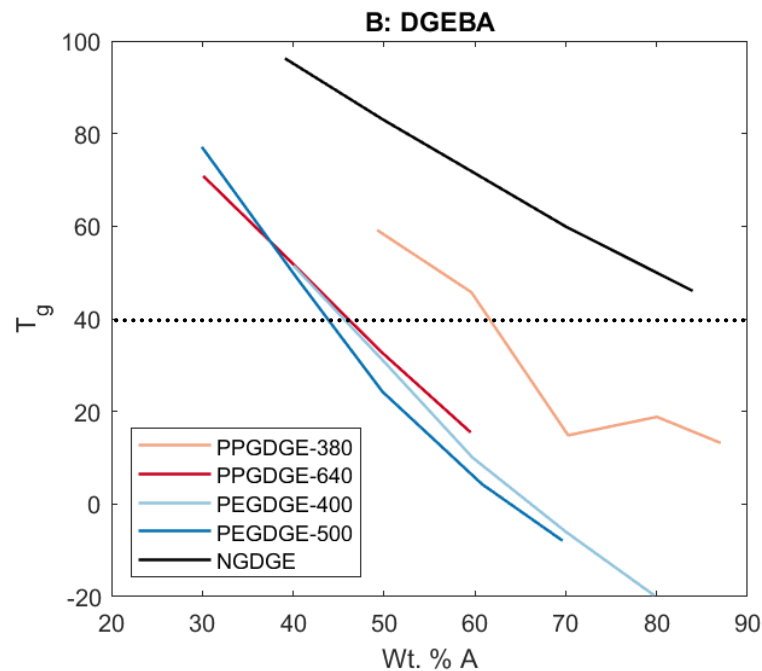
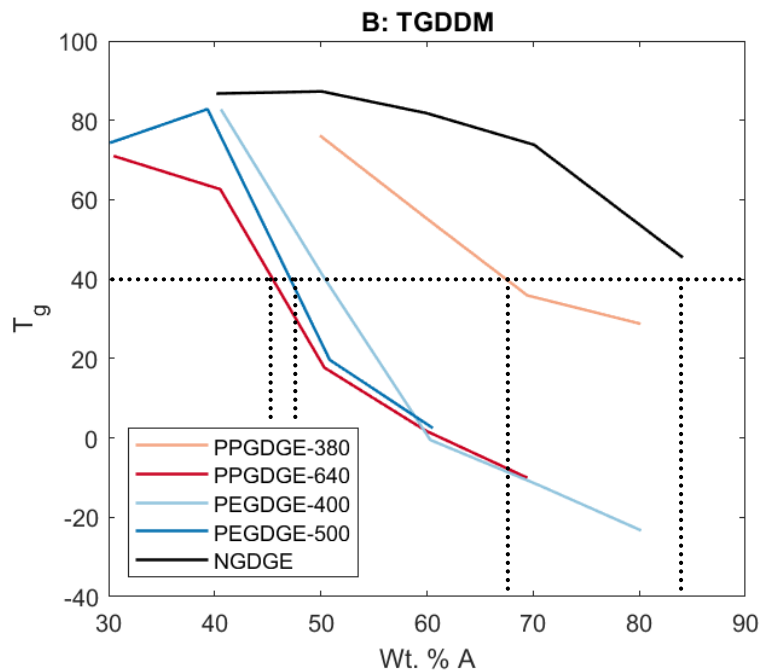
Systematic study – Results of prestudy

- Plots of T_g vs Wt. % component A
 - Fixed component A
 - Variable component B ($f = 2, 3, 4$)
- Findings
 - TGAP ($f=3$) seems to increase T_g the most, thus allowing to increase wt. % A
 - No advantage of TGDDM ($f=4$) over TGAP ($f=3$)
 - Effect not large



Systematic study – Results of prestudy

- Here, different molecular weights (Comp. A) are plotted together
- Findings
 - Higher molecular weight \rightarrow T_g decreases faster



Systematic study – Calculations

- Calculation of the crosslink density (under full conversion)

$$v_e = \frac{3}{2}C_3 + \frac{4}{2}C_4 + \frac{5}{2}C_5 + \dots = \sum_{f=3}^{\infty} C_f$$

C_f : Concentration of reactant of functionality f [moles/cm³]

v_e : Elastically active crosslinks [moles/cm³]

- Calculation of the corresponding rubbery elastic modulus (for small strains)

$$E = 3gv_eRT$$

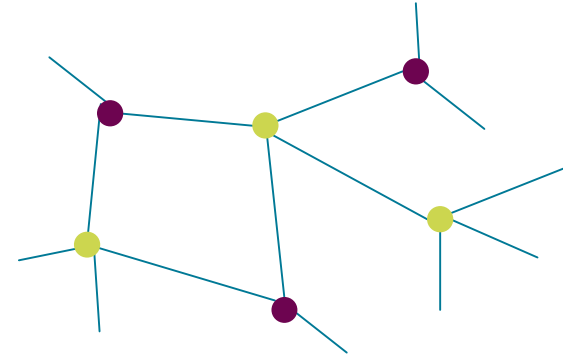
$$g = (f_c - 2)/f_c$$

g : Prefactor accounting for crosslink mobility

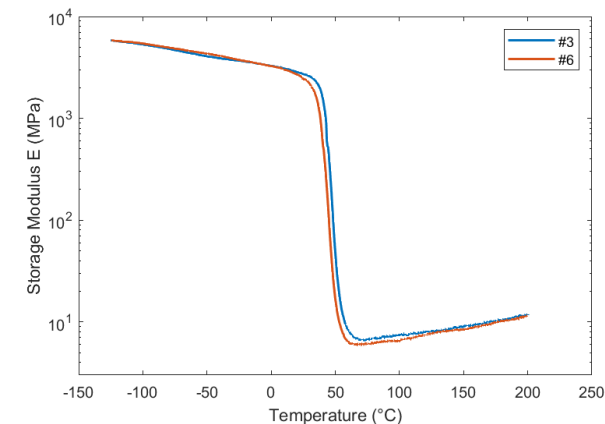
R : Gas constant

T : Temperature

f_c : Average crosslink functionality

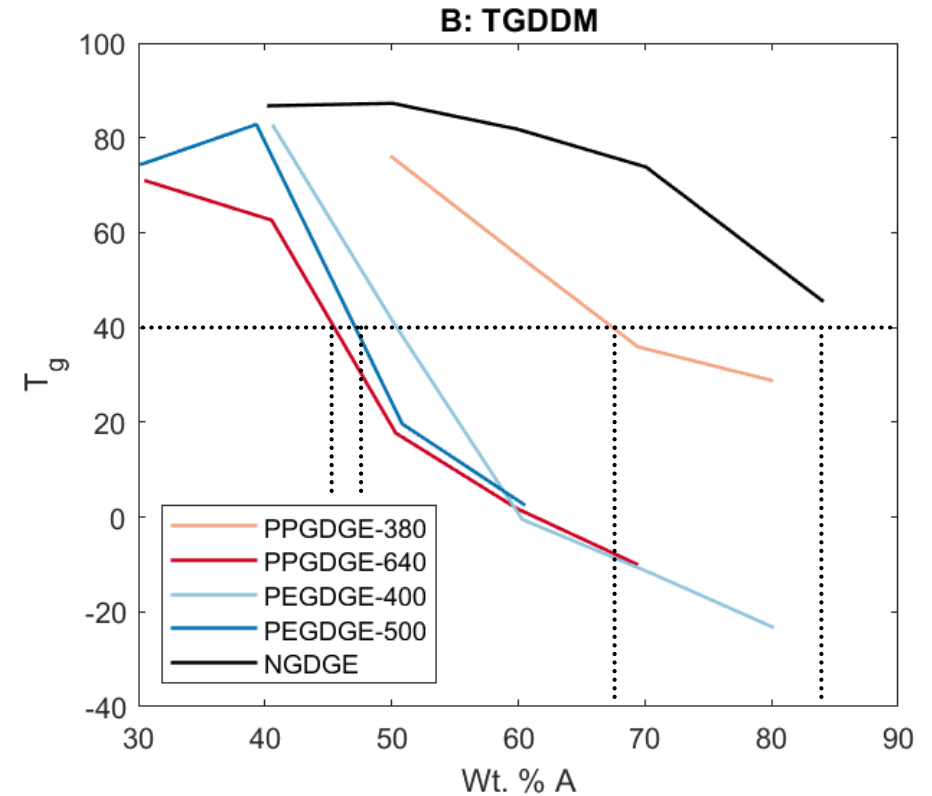


- Compare measured E , dE/dT with calculated value



Systematic study – Outlook

- Mechanical Testing: 27. Sept – 8. October
 - Toughness
 - DMTA
 - Compression
 - Tensile (above T_g)
- Investigation topological parameters - toughness
- Providing best formulations for PSI
 - All components are from Sigma-Aldrich



Acknowledgements

- ETH Zürich
 - Prof. Vermant, co-supervisor for discussions and support
 - Soft Materials Group for support
 - Xavier Guichard and Dr. Alessandro Lauria for synthesis and discussions on Gd_2O_3 nanoparticles
- PSI
 - Dr. Bernhard Auchmann, Dr. André Brem and Dr. Michael Daly for discussions and experimental support
- CERN
 - Dr. Davide Tomassini for discussions
- Funding by CHART (Swiss accelerator research and technology)
 - Collaboration between ETH, PSI, CERN, EPFL
 - Developing technology for the FCC (future circular collider) planned for 2060

