

Empa GEMS workshop 2014



Tutorial:

Influence of limestone on the hydration of calcium sulfoaluminate cements

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Thanks to: Barbara Lothenbach, Laure Pelletier-Chaignat & Lukas Martin

GEMS V.3.2 2482.930, CEMDATA14 (version April 14, 2014)

Influence of limestone on the hydration of calcium sulfoaluminate cements

1. Introduction
2. Hydrates in the system $C_4A_3S - Cs - H_2O$
3. Hydrates in the system $C_4A_3S - Cc - H_2O$
4. Hydrates in the system $C_4A_3S - Cs - Cc - H_2O$
5. Hydration products of a commercial CSA clinker in the presence of anhydrite and limestone
6. Appendix: occurrence of CAH_{10}

1. Introduction: CSA cements

Calcium sulfoaluminate cement (CSA)

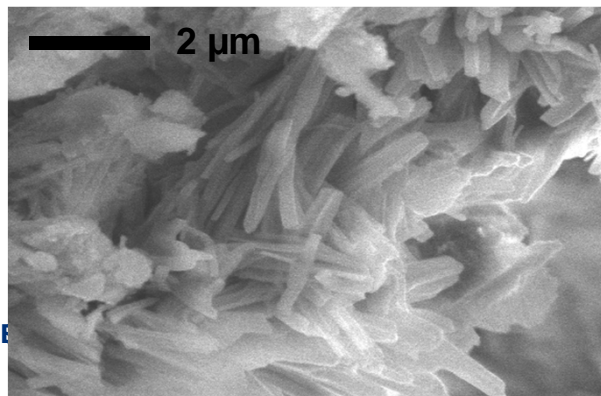
≈ 75-85% CSA clinker

- made from mixtures of calcite, clay and anhydrite (rotary kiln, ≈ 1250°C)
- main phase: ye'elimite = $4\text{CaO} \cdot 3\text{Al}_2\text{O}_3 \cdot \text{SO}_3$
- minor phases: belite, ferrite, calcium aluminates, perovskite, gehlenite, calcium sulfosilicate, ...

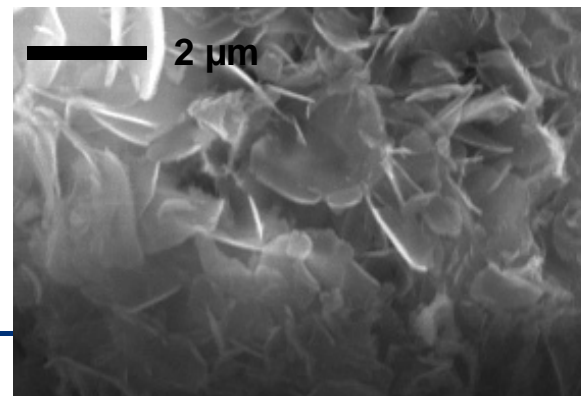
≈ 15-25% calcium sulfate (gypsum, anhydrite)

Main hydration products:

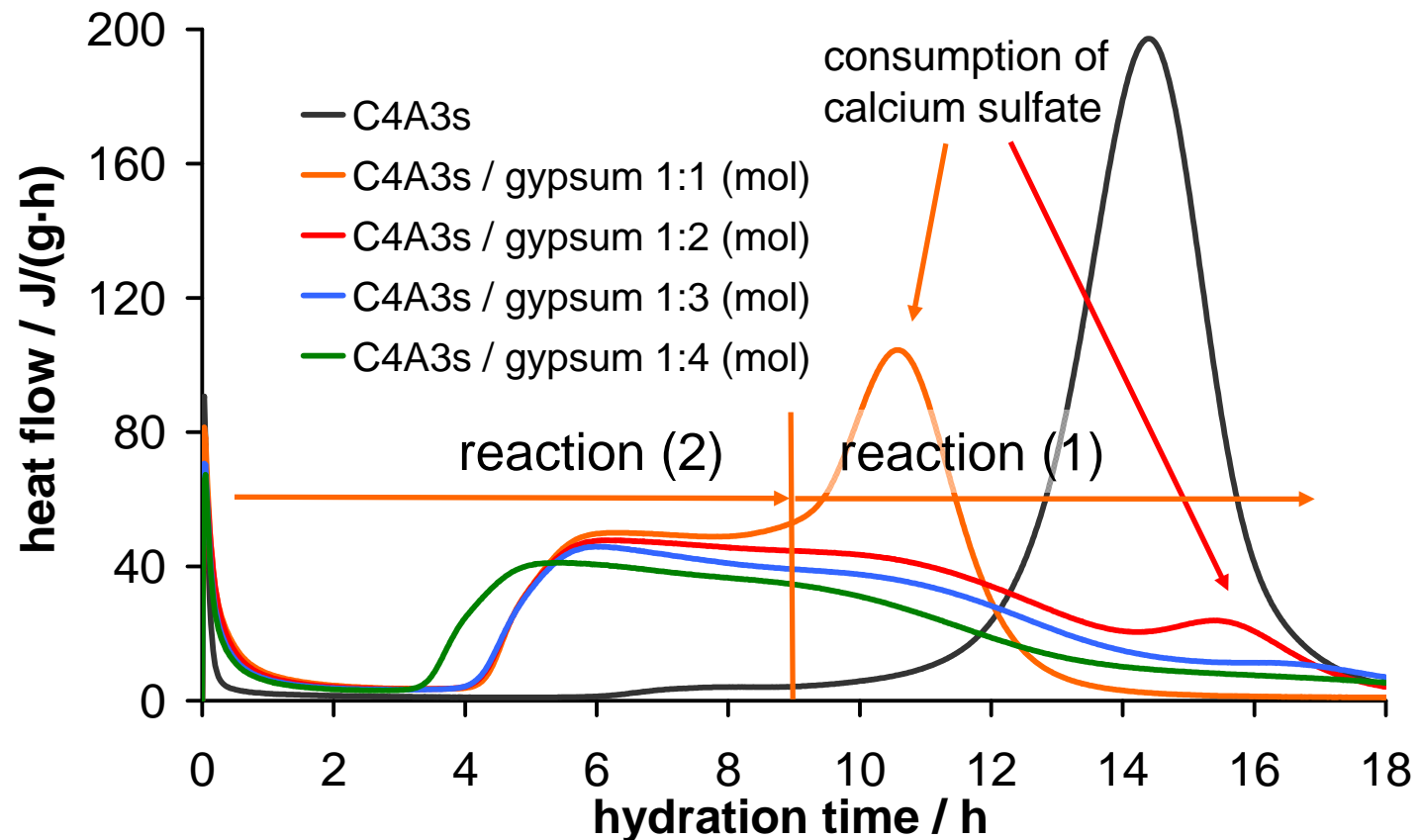
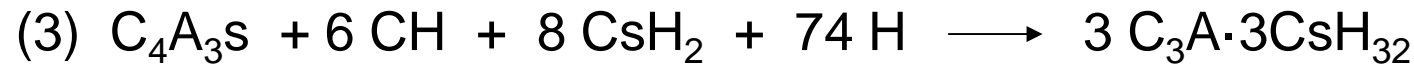
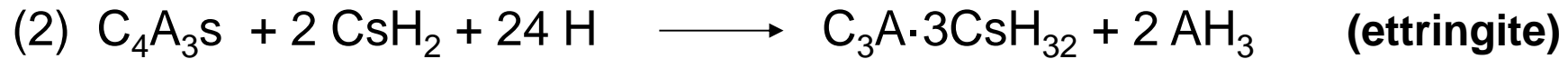
- ettringite,
 $3\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot 3\text{CaSO}_4 \cdot 32\text{H}_2\text{O}$



- monosulfate,
 $3\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot \text{CaSO}_4 \cdot 12\text{H}_2\text{O}$



1. Introduction: Hydration of ye'elimite



2. Hydrates in the system $C_4A_3s - Cs - H_2O$



Hydration of ye'elimite at 20°C

- variable amounts of **anhydrite**
- water/cement ratio of 1.00 (100 g solid and 100 g water)
- assumption of 100% hydration
- exclusion of CO_2
- single (parent file)
- process file for variable anhydrite additions
- comparison to experimental data

=> **open GEMS project "Yeelimite"**

2. Hydrates in the system $C_4A_3s - Cs - H_2O$



open single file „Ye_anh“:

hydration of a mixture of 90 g C_4A_3s + 10 g Cs + 100 g water + 1 g O_2

Input: System Definition		Results: Equilibrium State						
Phase/species	L	T	On/	UC	Add to BC	UG	G0 corr.	
aq_gen	29	a +	g	0		J	0	
gas_gen	5	g +	g	0		J	0	
SO4_OH_AFm	2	s +	g	0		J	0	
OH_SO4_AFm	2	s +	g	0		J	0	
SO4_CO3_AfT	2	s -	g	0		J	0	
CO3_SO4_AfT	2	s -	g	0		J	0	
Al(OH)3mic	1	s +	g	0		J	0	
Gibbsite	1	s -	g	0		J	0	
Graphite	1	s -	g	0		J	0	
Mayenite	1	s +	g	0		J	0	
Aluminate	1	s +	g	0		J	0	
CA	1	s +	g	0		J	0	
CA2	1	s +	g	0		J	0	
C2AH75	1	s +	g	0		J	0	
C3AH6	1	s +	g	0		J	0	
C4AH13	1	s -	g	0		J	0	
C4AH19	1	s +	g	0		J	0	
CAH10	1	s +	g	0		J	0	
C4AsH12	1	s -	g	0		J	0	
C4Ac0.5H12	1	s -	g	0		J	0	
C4AcH11	1	s -	g	0		J	0	
ettringite	1	s +	g	0		J	0	
Aragonite	1	s -	g	0		J	0	
Calcite	1	s -	g	0		J	0	
lime	1	s +	g	0		J	0	
Portlandite	1	s +	g	0		J	0	
Anhydrite	1	s +	g	0		J	0	
Gypsum	1	s +	g	0		J	0	
hemihydrate	1	s +	g	0		J	0	
Sulphur	1	s +	g	0		J	0	

← SO_4/CO_3 -AfT ss deactivated

← gibbsite deactivated
(does not form at ambient
temperature and „normal“
hydration times)

← we use the ss instead

all carbon containing phases
deactivated

2. Hydrates in the system $C_4A_3s - Cs - H_2O$



Result: hydration of a mixture of 90 g C_4A_3s + 10 g Cs + 100 g water + 1 g O_2

Input: System Definition		Results: Equilibrium State				
Phase/species	L	T	Amount (mol)	logSI/Activity	Concentration	
+ a aq_gen	23	a	2.1481191	1.75e-09		
+ g gas_gen	3	g	0.03119591	2.369e-10		
- s SO4_OH_AFm	2	s	0.10938991	3.174e-09		
C4AH13		J	0.0012490694	0.00134535	0.011418506	
monosulphate		M	0.10814084	0.987658	0.98858149	
- s OH_SO4_AFm	2	s	2.8306647e-007	3.174e-09		
C4AH13		M	3.2321935e-009	0.00134535	0.011418496	
monosulphate		J	2.7983428e-007	0.987658	0.9885815	
+ s Al(OH)3mic	1	s	0.59081956	-3.845e-10		
+ s Mayenite	1	s	0	-118.9		
+ s Aluminate	1	s	0	-37.14		
+ s CA	1	s	0	-8.171		
+ s CA2	1	s	0	-7.78		
+ s C2AH75	1	s	0	-0.7903		
+ s C3AH6	1	s	0	-0.8406		
+ s C4AH19	1	s	0	-2.413		
+ s CAH10	1	s	0	-0.2867		
+ s ettringite	1	s	0.037596495	-5.961e-09		
+ s lime	1	s	0	-11.49		
+ s Portlandite	1	s	0	-1.519		
+ s Anhydrite	1	s	0	-3.756		
+ s Gypsum	1	s	0	-3.482		
+ s hemihydrate	1	s	0	-4.529		
+ s Sulphur	1	s	0	-120.5		

solid solution
 SO_4/OH -AFm

microcrystalline
 AH_3

ettringite

=> we will now look at the process file

2. Hydrates in the system $C_4A_3s - Cs - H_2O$



open process file „Ye_anh_pr“:

hydration of a mixture of

100 g ($C_4A_3s + Cs$; various C_4A_3s/Cs ratios) + 100 g water + 1 g O_2

Input

Controls | Sampling | Results | Config | 21/04/2014, 15:51

Addition of anhydrite to ye'elimite
GEMS 2014

101 single systems

	iTm	iV	iP	iTC	iNv	iTau	ipXi	iNu	ipH	ipe
0	1000	0	1	20	0	0	0	0	0	0
1	1100	0	1	20	0	0	0	50	0	0
2	1	0	0	0	0	0	0	0.5	0	0
cTm	1100	0	1	20	0	0	0	50	0	0

```
$ hydration of 100 g solid with 100 g water and 1 g oxygen added
xa_{Aqua} =: 100;
xa_{O2} =: 1;
$ amount of calcium sulfate is taken from iNu
xa_{CaSO4} =: cNu;
$
$ calculation of the amount of ye'elimite
xa_{C4A3s} =: 100-cNu;
```

creates numbers 0 ... 50 in steps of 0.5
= % Cs in the system)

2. Hydrates in the system $C_4A_3s - Cs - H_2O$



Sampling

Controls	Sampling	Results	Config	21/04/2014, 15:51							
NeIt	9999	101	Next	0	I	0	J	100	Jp	100	
pSTkey				Yeelimite:G:Ye_anh:0:0:1:20:0:				cTm	1100	cNV	0
cTau	0	cpXi	0	cXi	1	cNu	50				
cpH	0	cpe	0	cEh	0	cT	293.15				

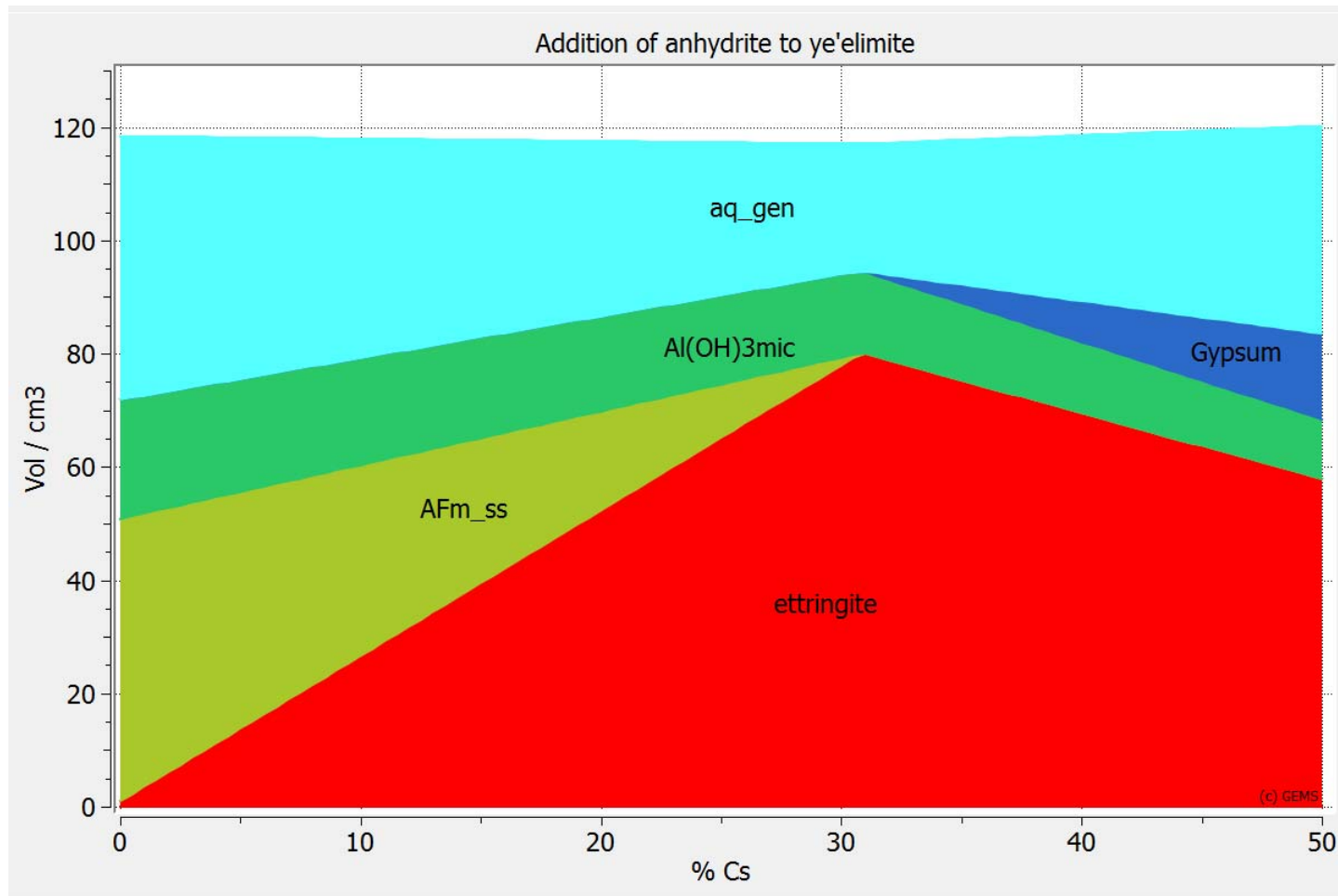
```
$ x-axis: amount of anhydrite added
xp[J] =: cNu;
$ y-axis: phase volumes
yp[J][0] =: phVol[{ettringite}];
yp[J][1] =: phVol[{SO4_OH_AFm}]+phVol[{OH_SO4_AFm}];
yp[J][2] =: phVol[{Al(OH)3mic}];
yp[J][3] =: phVol[{Gypsum}];
yp[J][4] =: phVol[{aq_gen}];
```

we need both solid solutions

=> calculate systems using the process file and display results graph

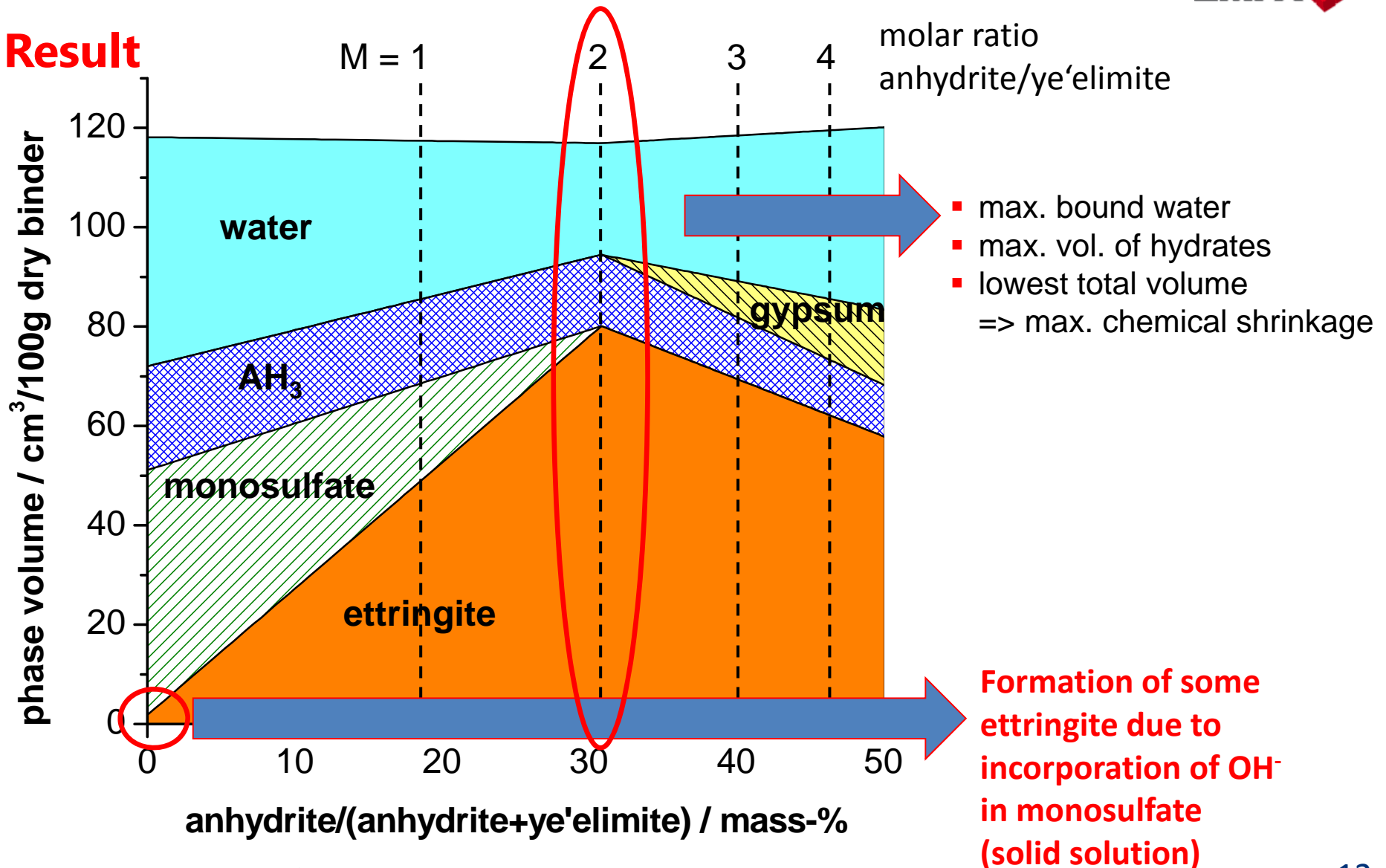
2. Hydrates in the system $C_4A_3s - Cs - H_2O$

Result



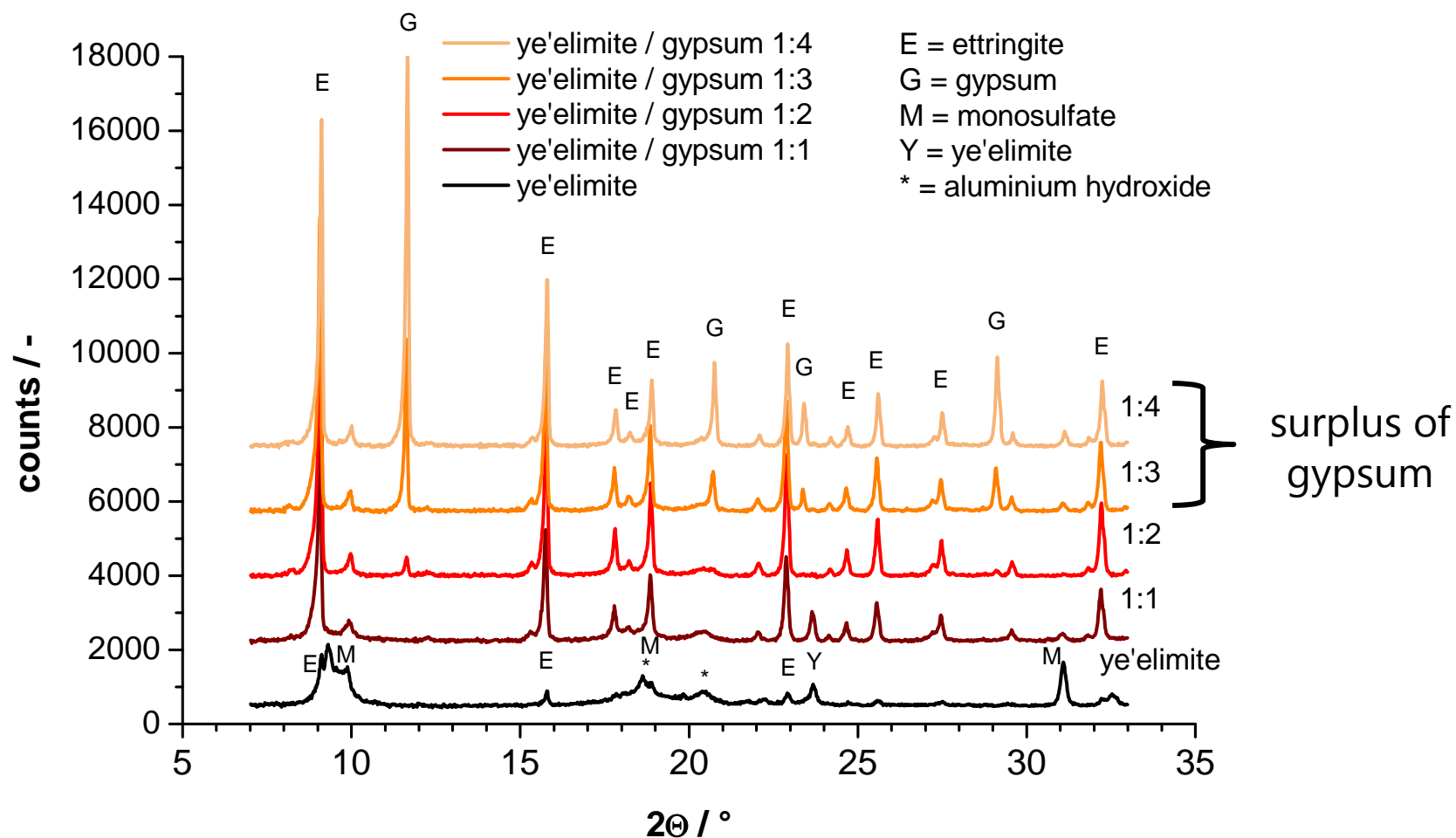
2. Hydrates in the system $C_4A_3s - Cs - H_2O$

Result



2. Hydrates in the system $C_4A_3S - Cs - H_2O$

Experiment: XRD after 18 h of hydration at water/solid = 2



3. Hydrates in the system $C_4A_3S - Cc - H_2O$



Hydration of ye'elimite at 20°C

- variable amounts of **calcite**
- water/cement ratio of 1.00 (100 g solid and 100 g water)
- assumption of 100% hydration
- exclusion of CO_2
- single (parent file)
- process file for variable calcite additions
- comparison to experimental data

=> **open single file "Ye_cal" in GEMS project "Yeelimite"**

3. Hydrates in the system $C_4A_3s - Cc - H_2O$



open single file „Ye_cal“:

hydration of a mixture of 90 g C_4A_3s + 10 g Cc + 100 g water + 1 g O_2

Input: System Definition					Results: Equilibrium State				
Phase/species	L	T	On/	UC					
aq_gen	29	a	+	g					
gas_gen	5	g	+	g					
SO4_OH_AFm	2	s	+	g					
OH_SO4_AFm	2	s	+	g					
SO4_CO3_Aft	2	s	+	g					
CO3_SO4_Aft	2	s	+	g					
Al(OH)3mic	1	s	+	g					
Gibbsite	1	s	-	g					
Graphite	1	s	+	g					
Mayenite	1	s	+	g					
Aluminate	1	s	+	g					
CA	1	s	+	g					
CA2	1	s	+	g					
C2AH75	1	s	+	g					
C3AH6	1	s	+	g					
C4AH13	1	s	-	g					
C4AH19	1	s	+	g					
CAH10	1	s	+	g					
C4AsH12	1	s	-	g					
C4Ac0.5H12	1	s	+	g					
C4AcH11	1	s	+	g					
ettringite	1	s	-	g					
Aragonite	1	s	+	g					
Calcite	1	s	+	g					
lime	1	s	+	g					
Portlandite	1	s	+	g					
Anhydrite	1	s	+	g					
Gypsum	1	s	+	g					
hemihydrate	1	s	+	g					
Sulphur	1	s	+	g					

deactivated
(metastability
constraint)

deactivated
(use of ss
instead)

Compos	DComp	Phase	IComp	Surfaces	Config	21/04/2014, 12:
Hydration of ye'elimite + calcite						
GEMS 2014						
Masses	0	1	1	1	0.201	21.1965
Volumes	1	1	Procs	0	0	0
AC	CCvp				xa	
0	*	Al(OH)3	MIN Aluminum-hydroxide...	g		0
1	*	Al2O3	MIN Aluminum-oxide_1M_	g		0
2	+	Aqua	AQ 1_mole_H2O_	g		100
3	*	C12A7	MIN Mayenite_1M_	g		0
4	*	C3A	MIN Tricalcium_alumina...	g		0
5	+	C4A3s	MIN Yeelimite_1M_	g		90
6	*	CA	MIN Calcium_aluminate_...	g		0
7	*	CA2	MIN Calcium_dialuminat...	g		0
8	*	CH4	GA Methane_1M_	g		0
9	*	CO2	GA Carbon-dioxide_1M_	g		0
10	*	Ca(OH)2	MIN Calcium-hydroxide_...	g		0
11	+	CaCO3	MIN Calcium-carbonate_...	g		10
12	*	CaO	MIN Calcium-oxide_1M_	g		0
13	*	CaSO4	MIN Calcium-sulfate_1M_	g		0
14	*	CaSO4_05H2O	MIN hemihydrate_1M_	g		0
15	*	Gypsum	MIN Ca-sulfate-2H2O-1M_	g		0
16	*	H2	GA Hydrogen_1M_	g		0
17	*	H2S	GA Hydrogen-sulfide_1...	g		0
18	*	H2SO4	AQ Sulfuric-acid_1M_	g		0
19	+	O2	GA Oxygen_1M_	g		1
20	*	SO3	MIN Sulfur-trioxide_1M_	g		0

=> calculate phase assemblage

3. Hydrates in the system $C_4A_3s - Cc - H_2O$



Result

Input: System Definition		Results: Equilibrium State				
Phase/species	L	T	Amount (mol)	logSI/Activity	Concentration	
+ a aq_gen	29	a	1.9949533	9.829e-10		
+ g gas_gen	5	g	0.031199785	-1.343e-09		
+ s SO4_OH_AFm	2	s	0	-0.6997		
+ s OH_SO4_AFm	2	s	0	-0.6997		
- s SO4_CO3_AfT	2	s	0.049981584	3.211e-08		
+ tricarboalu		J	0.00082269489	0.000580702	0.01645996	
+ ettringite		M	0.049158889	0.981119	0.98354004	
- s CO3_SO4_AfT	2	s	4.1696143e-007	3.211e-08		
+ tricarboalu		M	6.8631924e-009	0.000580702	0.016460018	
+ ettringite		J	4.1009824e-007	0.981119	0.98353998	
+ s Al(OH)3mic	1	s	0.58996332	5.526e-10		
+ s Graphite	1	s	0	-82.72		
+ s Mayenite	1	s	0	-123.1		
+ s Aluminate	1	s	0	-38.18		
+ s CA	1	s	0	-8.517		
+ s CA2	1	s	0	-8.126		
+ s C2AH75	1	s	0	-1.483		
+ s C3AH6	1	s	0	-1.88		
+ s C4AH13	1	s	0	-4.256		
+ s C4AH19	1	s	0	-3.798		
+ s CAH10	1	s	0	-0.6328		
+ s C4AsH12	1	s	0	-0.7009		
+ s C4Ac0.5H12	1	s	0	-1.248		
+ s C4AcH11	1	s	0.097444931	-5.071e-09		

solid solution
SO₄/CO₃-AfT

microcrystalline
AH₃

monocarbonate

=> we will now look at the process file

3. Hydrates in the system $C_4A_3s - Cc - H_2O$



open process file „Ye_cal_pr“:

hydration of a mixture of 100 g ($C_4A_3s + Cc$; various C_4A_3s/Cc ratios)

+ 100 g water + 1 g O_2

similar structure than previous file

Input

```
$ hydration of 100 g solid with 100 g water and 1 g oxygen added
xa_[{Aqua}] =: 100;
xa_[{O2}] =: 1;
$ amount of calcium carbonate is taken from iNu
xa_[{CaCO3}] =: cNu;
$
$ calculation of the amount of ye'elimite
xa_[{C4A3s}] =: 100-cNu;
```

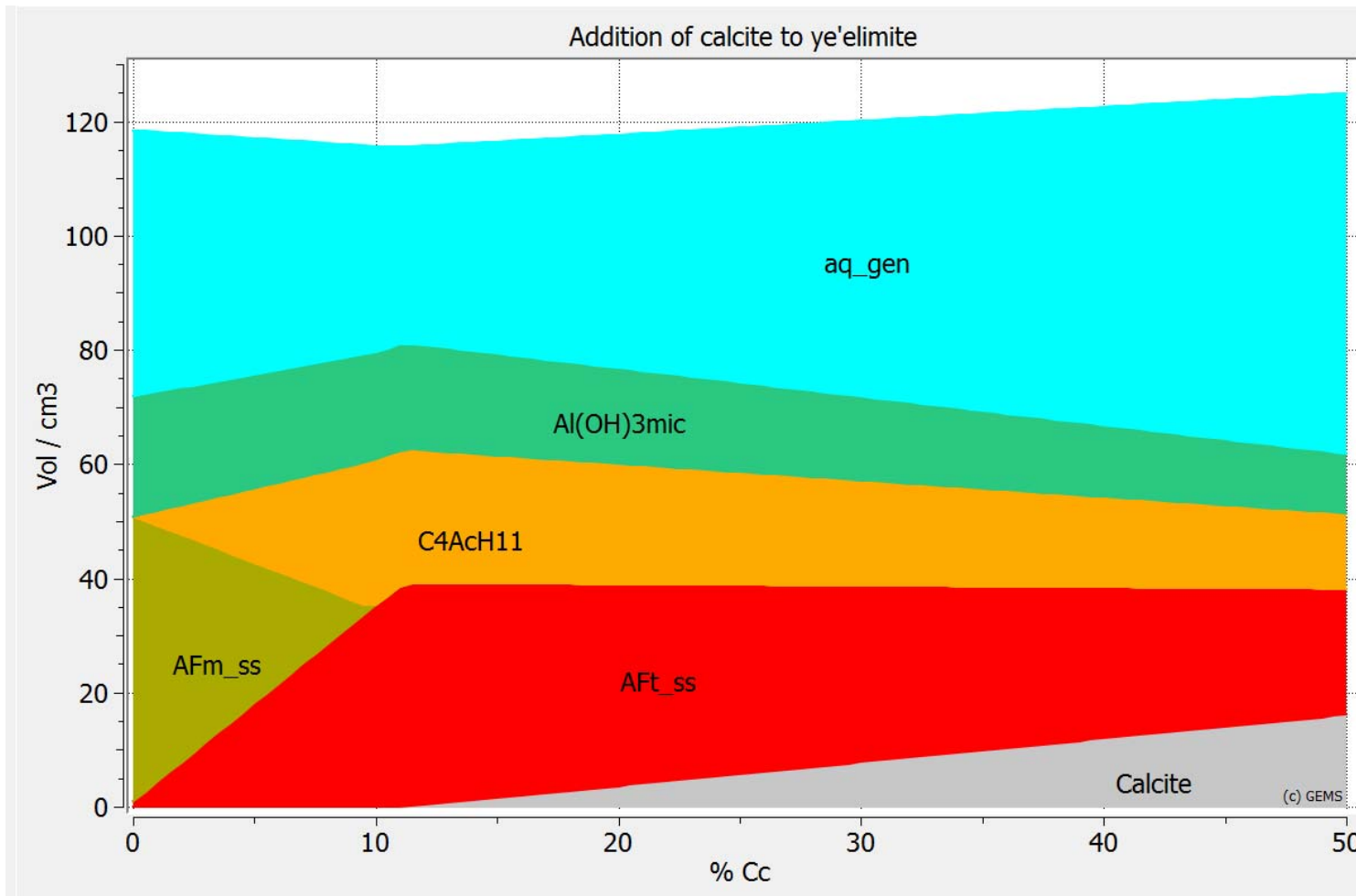
Output

```
$ x-axis: amount of calcite added
xp[J] =: cNu;
$ y-axis: phase volumes
yp[J][0] =: phVol[{Calcite}];
yp[J][1] =: phVol[{SO4_CO3_AfT}]+phVol[{CO3_SO4_AfT}];
yp[J][2] =: phVol[{SO4_OH_AfM}]+phVol[{OH_SO4_AfM}];
yp[J][3] =: phVol[{C4AcH11}];
yp[J][4] =: phVol[{Al(OH)3mic}];
yp[J][5] =: phVol[{aq_gen}];
```

=> calculate systems using the process file and display results graph

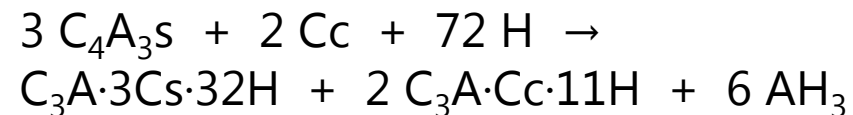
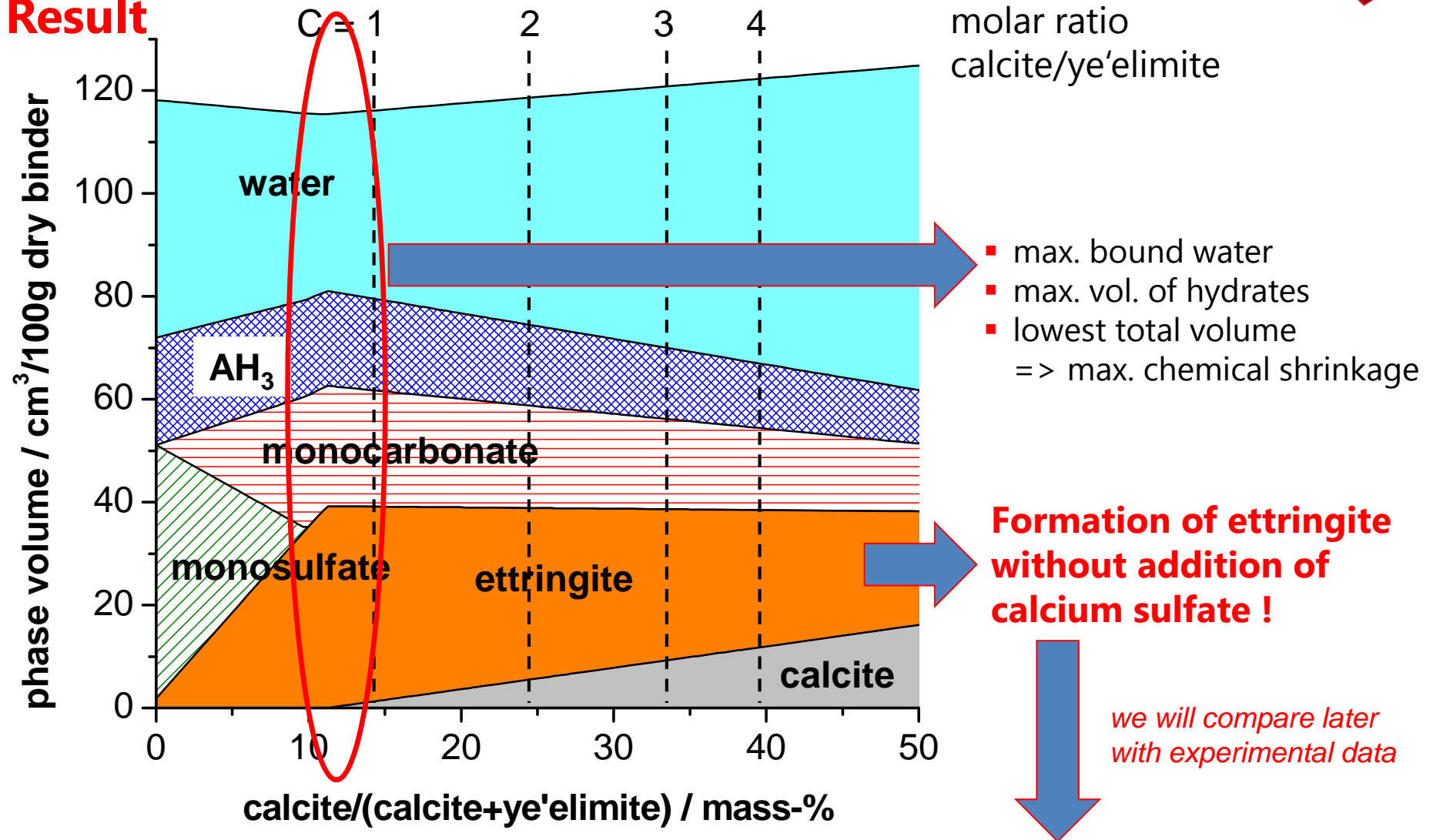
3. Hydrates in the system $C_4A_3s - Cc - H_2O$

Result



3. Hydrates in the system $C_4A_3s - Cc - H_2O$

Result



4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$



Hydration of ye'elimite at 20°C

- variable amounts of **anhydrite and calcite**
- water/cement ratio of 1.00 (100 g solid and 100 g water)
- assumption of 100% hydration
- exclusion of CO_2
- 2 process files for variable calcite or anhydrite additions
- comparison to experimental data

=> **open single file "Ye_anh_cal_1" in GEMS project**
"Yeelimite": constant calcite, varying anhydrite
(Ye_anh_cal_2: vice versa with constant calcite, not further discussed in this presentation)

4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$



open process file „Ye_anh_cal_1“:

hydration of a mixture of 100 g (86 g $C_4A_3s + Cs$ + 14 g Cc ;
varying C_4A_3s/Cs ratios) + 100 g water + 1 g O_2
(process file *Ye_anh_cal_2* does the same vice versa with fixed Cs content)

Input

```
$ hydration of 100 g solid with 100 g water and 1 g oxygen added
xa_{{Aqua}} =: 100;
xa_{{O2}} =: 1;
$
$ amount of calcium carbonate in g to be changed below
xa_{{CaCO3}} =: 14;
$
$ amount of calcium sulfate is taken from iNu
xa_{{CaSO4}} =: cNu;
$
$ calculation of the amount of ye'elimite
xa_{{C4A3s}} =: 100-xa_{{CaCO3}}-cNu;
```

change here if amount of calcite
should be varied

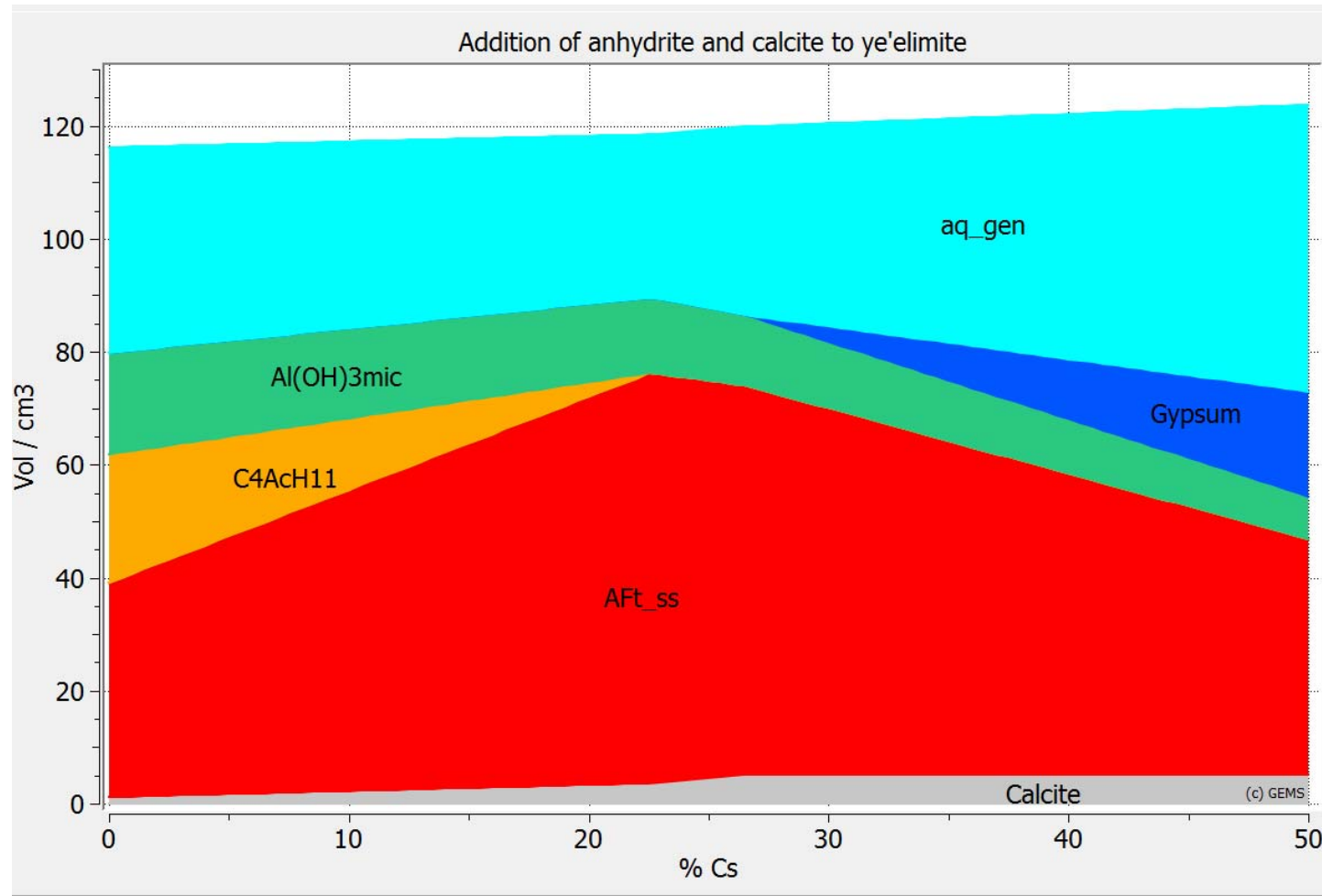
Output

```
$ x-axis: amount of anhydrite added
xp[J] =: cNu;
$ y-axis: phase volumes
yp[J][0] =: phVol[{{Calcite}}];
yp[J][1] =: phVol[{{SO4_CO3_Aft}}]+phVol[{{CO3_SO4_Aft}}];
yp[J][2] =: phVol[{{SO4_OH_AfM}}]+phVol[{{OH_SO4_AfM}}];
yp[J][3] =: phVol[{{C4ACH11}}];
yp[J][4] =: phVol[{{Al(OH)3mic}}];
yp[J][5] =: phVol[{{Gypsum}}];
yp[J][6] =: phVol[{{aq_gen}}];
```

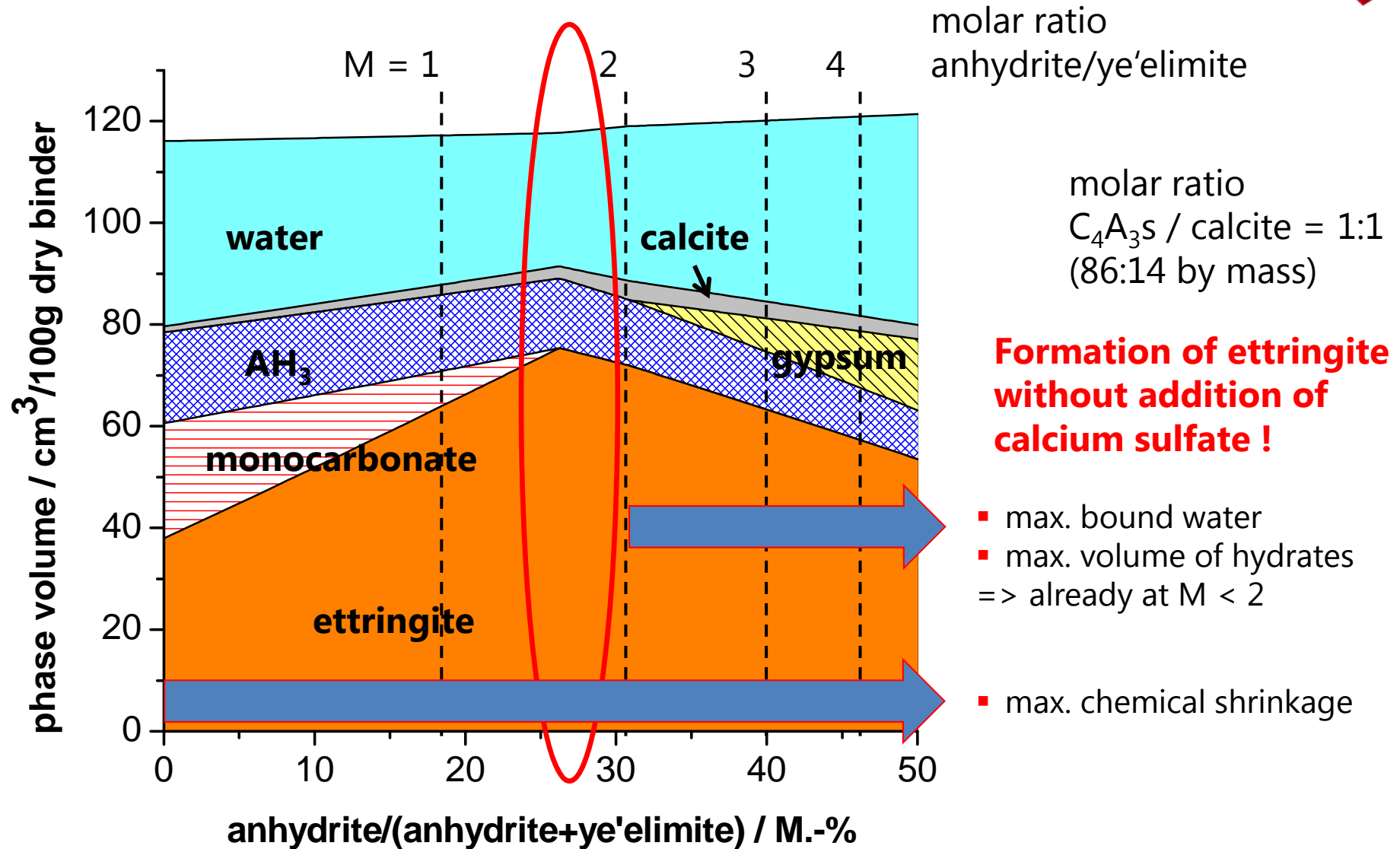
=> calculate systems using the process file and display results graph

4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$

Result



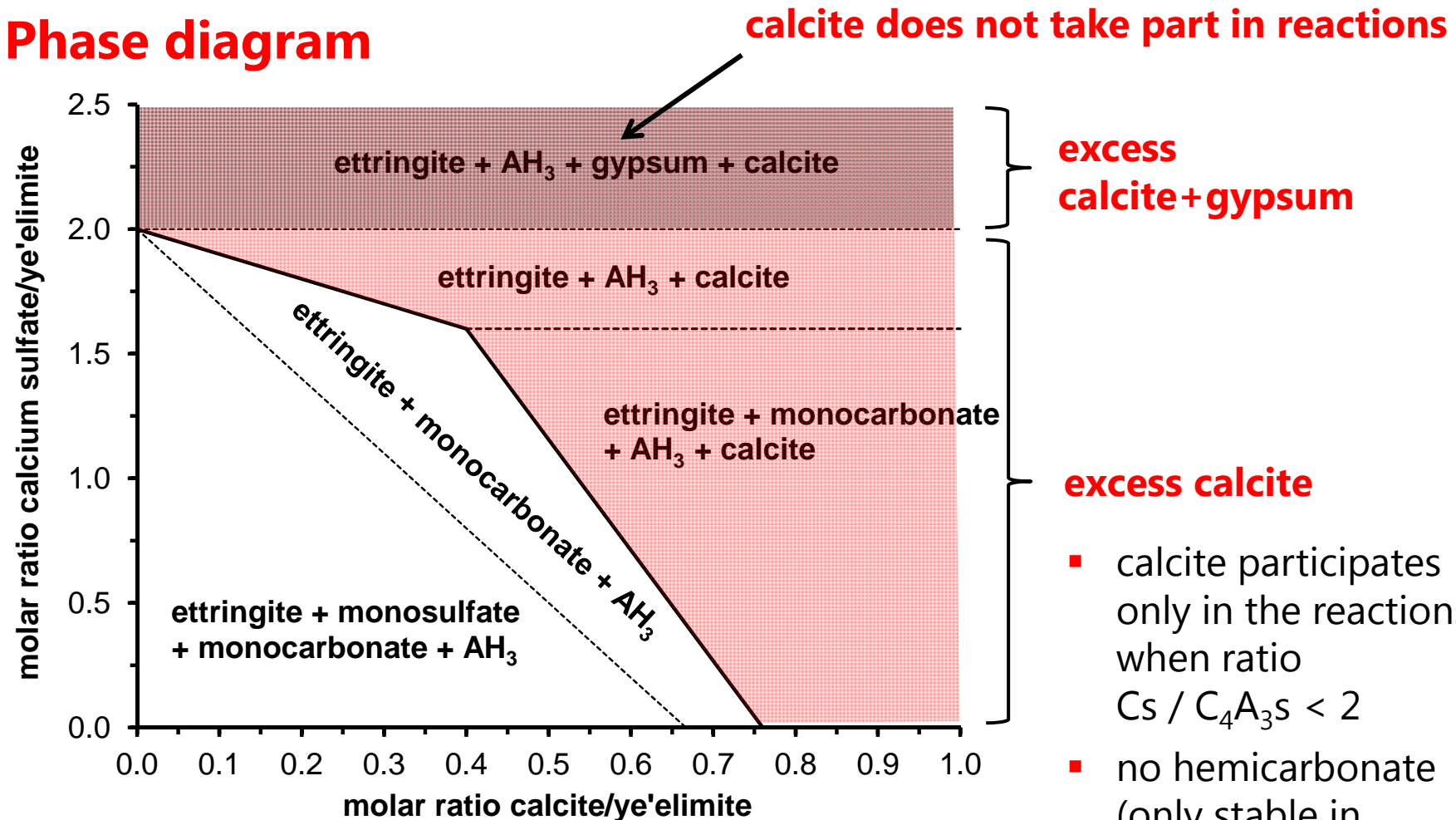
4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$



4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$



Phase diagram

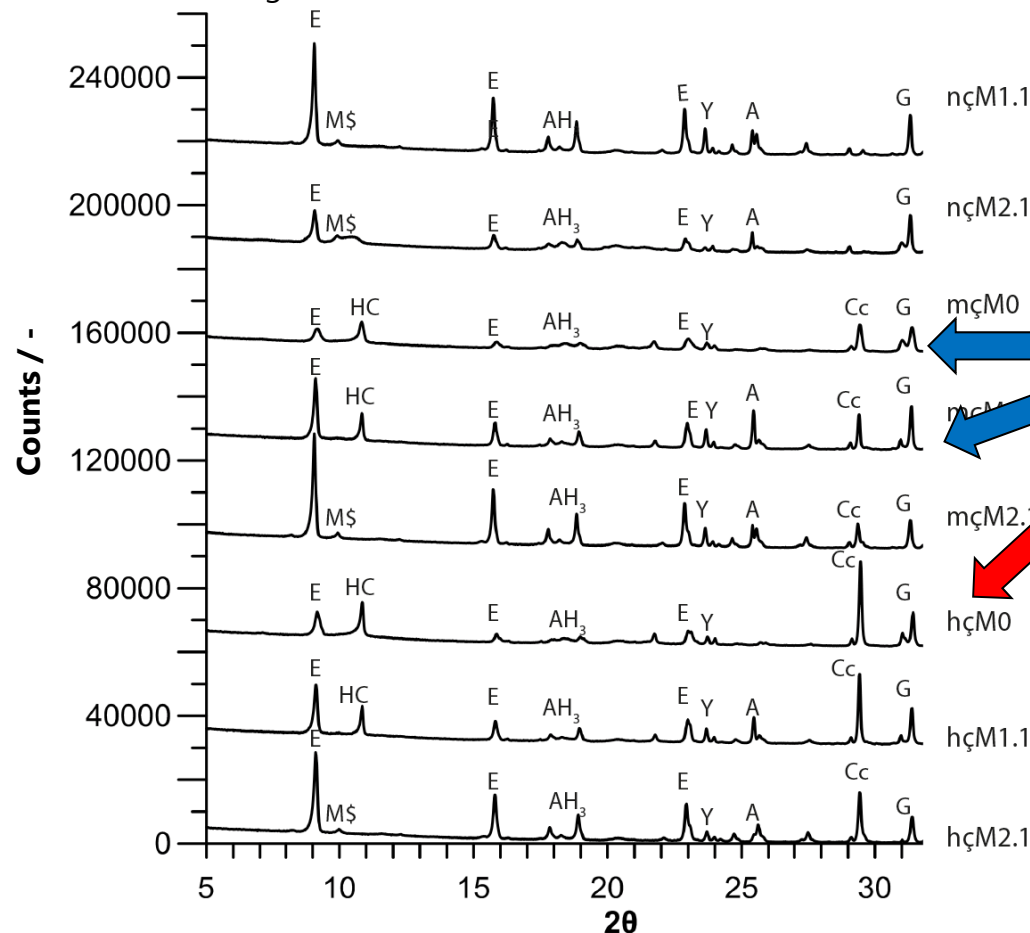


- calcite participates only in the reactions, when ratio $Cs / C_4A_3s < 2$
- no hemicarbonates (only stable in presence of portlandite => OPC)

4. Hydrates in the system $C_4A_3s - Cs - Cc - H_2O$

Experiment: XRD after 90 d of hydration (commercial CSA cement)

E = ettringite; AH_3 = aluminium hydroxide; Y = ye'elimite; A = anhydrite;
Cc = calcite; G = gehlenite;; MS = monosulfate; HC = hemicarbonate



Binder composition

Anhydrite Calcite

Low	High
-	-
High	-

- Participation of calcite **without anhydrite** (formation of ettringite) or when **low&medium amounts of anhydrite** are added

- Formation of hemicarbonate instead of monocarbonate (metastability)

Low	High
High	high

Martin L.H.J, Winnefeld F., Müller C.J., Lothenbach B.: 1st International Conference on Sulphoaluminate Cement: Materials and Engineering Technology, Wuhan, China, October 23-24, 2013, 229.

5. Commercial CSA clinker + anhydrite + limestone



Hydration of a commercial CSA clinker at 20°C

- variable amounts of anhydrite and calcite
- water/cement ratio of 1.00 (100 g solid and 100 g water)
- assumption of 100% hydration
- exclusion of CO₂
- 2 process files for variable calcite or anhydrite additions
- reactive part of the CSA clinker as predefined composition in GEMS

5. Commercial CSA clinker + anhydrite + limestone



Composition of the CSA clinker

needs to be recalculated
(reactive part only) in mol per
100 g referred to the elements
rough assumption:
inert phases contain no minor
elements

XRF		QXRD	
	CSA clinker		CSA clinker
SiO ₂	5.41	C ₄ A ₃ S̄	68.1
Al ₂ O ₃	42.61	CA	3.2
Fe ₂ O ₃	1.55	CA ₂	0.7
Cr ₂ O ₃	0.032	C ₁₂ A ₇	1.4
MnO	0.023	C ₂ AS	19.4
TiO ₂	2.100	C ₂ S	1.7
P ₂ O ₅	0.113	CT	3.9
CaO	35.87	M	0.6
MgO	1.01	MA	1.1
K ₂ O	0.53		
Na ₂ O	0.03		
SO ₃	8.52		
L.O.I.	1.98		
Total	99.75		
Density (g/cm³)	2.75		
Blaine (cm²/g)	5080		
CaO_{free} (wt%)	0.06		

inert:
24.4%

- [1] Martin L.H.J., Winnefeld F., Müller C.J., Lothenbach B.: 1st International Conference on Sulphoaluminate Cement: Materials and Engineering Technology, Wuhan, China, October 23-24, 2013, 229.

5. Commercial CSA clinker + anhydrite + limestone



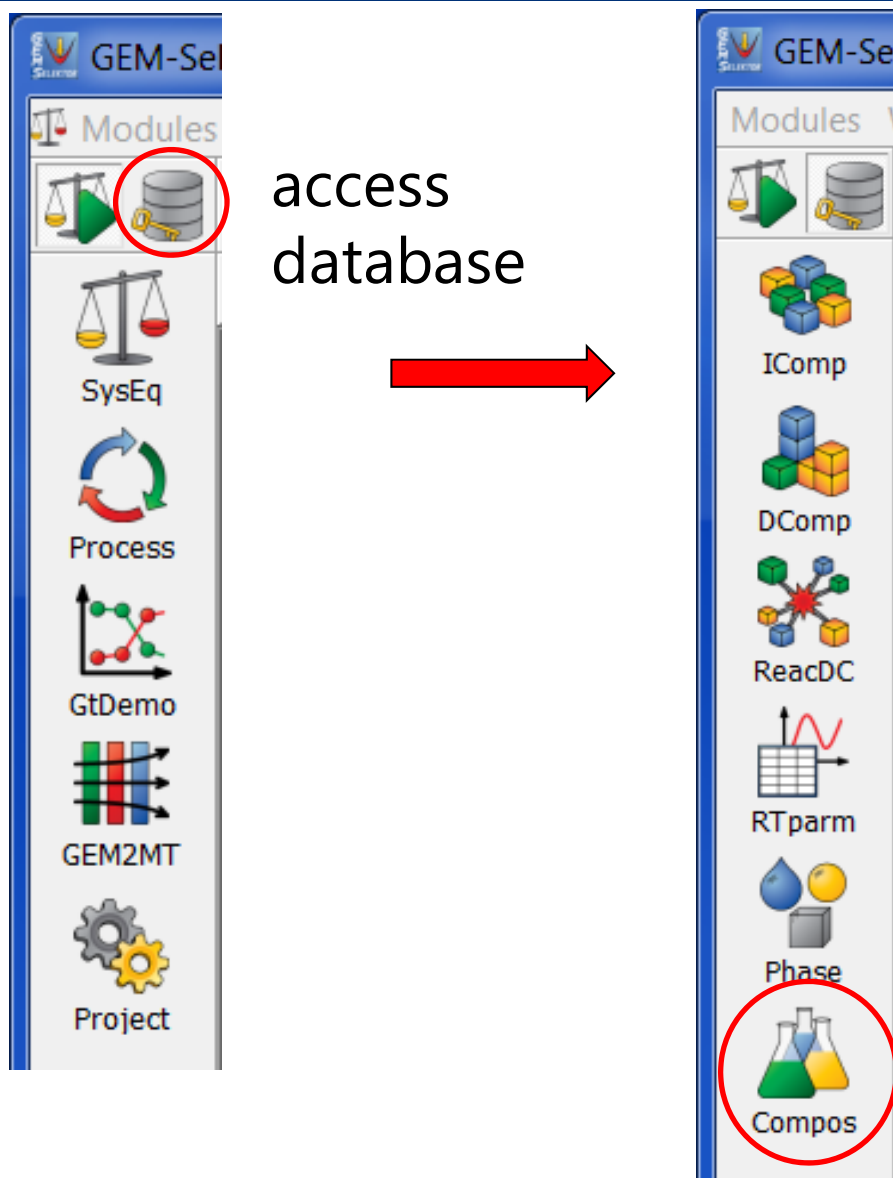
Calculation of the composition of the reactive part

Predefined composition									
CSA				nonreactive	reactive	normalized	Elements		
	XRD / M.-%		XRF / M.-%	M.-%	M.-%	M.-%		g/100g	mol/100g
C4A3s	68.10	SiO ₂	5.41	4.25	1.16	1.58263034	Si	0.73977669	0.02634016
CA	3.20	Al ₂ O ₃	42.61	8.00	34.61	47.2587898	Al	25.0117477	0.92699485
CA ₂	0.70	Fe ₂ O ₃	1.55	0.19	1.36	1.85591639	Fe	1.29808923	0.02324367
C ₂ AS *	19.40	CaO	35.87	9.54	26.33	35.9497673	Ca	25.6929668	0.64107408
CT * **	3.90	Na ₂ O	0.03		0.03	0.04096665	Na	0.03039141	0.00132195
MA *	1.10	K ₂ O	0.53		0.53	0.72374413	K	0.60081458	0.01536677
M	0.60	MgO	1.01	0.31	0.70	0.95365699	Mg	0.5750894	0.02366136
C ₁₂ A ₇	1.40	SO ₃	8.52		8.52	11.6345283	S	4.65967044	0.14531499
C ₂ S	1.70	TiO ₂	2.1	2.10	0.00	0	Ti	0	0
		H ₂ O			0.00	0	H	0	0
							O	41.3914538	2.58706287

75.6% is reactive, 24.4% is inert.

=> open now GEMS project "CSA"

5. Commercial CSA clinker + anhydrite + limestone



5. Commercial CSA clinker + anhydrite + limestone



GEM-Selektor 3 (GEMS3) - Geochemical Equilibrium Modelling by Gibbs Energy Minimization

Modules Record Record List Database Files Window Help

+ Create(New)...
New(Clone)...
Display F6
Remake...
Calculate F9
Save
Save As...
Delete
Plot data
Print...

IComp
DComp
ReacDC
RTparm
Phase
Compos

Al(OH)₃:MIN:Aluminum-hydroxide

Compos :: Predefined composition objects (PCO)

PCO Settings 22/04/2014, 09:56

Aluminum hydroxide Al(OH)₃ 1 mol
GEMS PCO database

0.0780036 0 0 0 0 0

	symIC	PCO		symIC	CIc	CI
0	Al ...	1	0	Al ...	M	1
1	H ...	3	1	H ...	M	3
2	O ...	3	2	O ...	M	3

5. Commercial CSA clinker + anhydrite + limestone



Compos: Please, set a new record key

CSA:MIN:CSA_clinker_1M_:

CSA Name of predefined composition object (PCO)

MIN Code of PCO type { AQ RO GA FL HC PM MIN }

CSA_clinker_1M_ Comment to PCO description

Ok Reset From List Help Cancel

5. Commercial CSA clinker + anhydrite + limestone



GEM-Selektor Compos Setup: CSA:MIN:CSA_clinker_1M_

Step 1 - Predefined Composition Object (PCO) configuration

Compos record contains data for a Predefined Composition Object (PCO), which describes a salt, the air, a mineral, a rock, a natural water, etc., treated as a single named entity in the chemical system recipe.

PCO can be configured according to the source data: as IC or DC amounts/concentrations; as user-defined formula (UDF) units; or as a large UDF entered into a text field.

PCO input data configuration

☒ Use amounts of Independent Components (IComp) in this PCO definition (default)

☐ Use formulae of Dependent Components (from DComp/ReacDC records) in this I

0 Set number of user-defined formula units for this PCO definition (0 by default)

Optional: Input user-defined formula (UDF) text

☐ Use a user-defined formula text input field for this PCO definition?

M moles Select units of measurement for this UDF quantity (default: M)

0 Enter here the UDF quantity or amount in selected units (default:

Learn more < Back **Next>** Cancel

GEM-Selektor Compos Setup: CSA:MIN:CSA_clinker_1M_

Step 2 - Additional settings and next actions

Optional

0 Set here the number of links to SDref bibliography records (default 0)

☐ Use a vector of data uncertainties in this PCO definition

What will happen after you click "Finish"

(1) For a PCO definition using IComp amounts, a list of available IComp records will be shown, asking you to mark the desired ones.

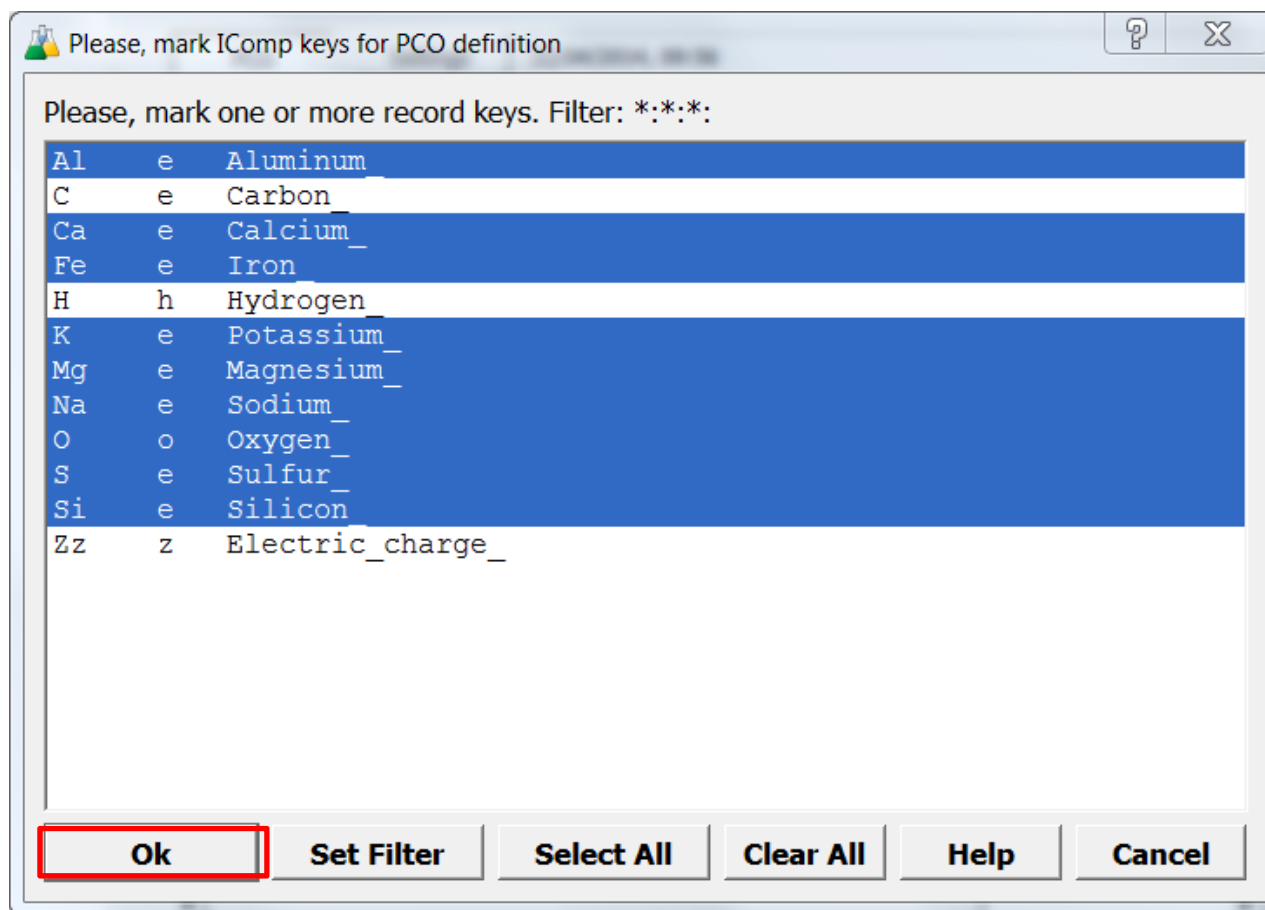
(2) In a PCO definition using DC formulae, a list of available ReacDC/DComp records will be shown, asking you to mark the desired ones.

(3) Page 1 of the 'Compos' window appears. Fill out BCname field and (optionally) BCnote lines. Then enter data and formulae wherever needed, check units of amount/concentration.

(4) Click on 'Calculate' toolbar button to compute the PCO vector. Check or enter the normalization values in MasVol[0] and MasVol[6] fields and calculate again, if needed. Setting both fields to zero disables the normalization of PCO to the total mass or total amount of moles, respectively.

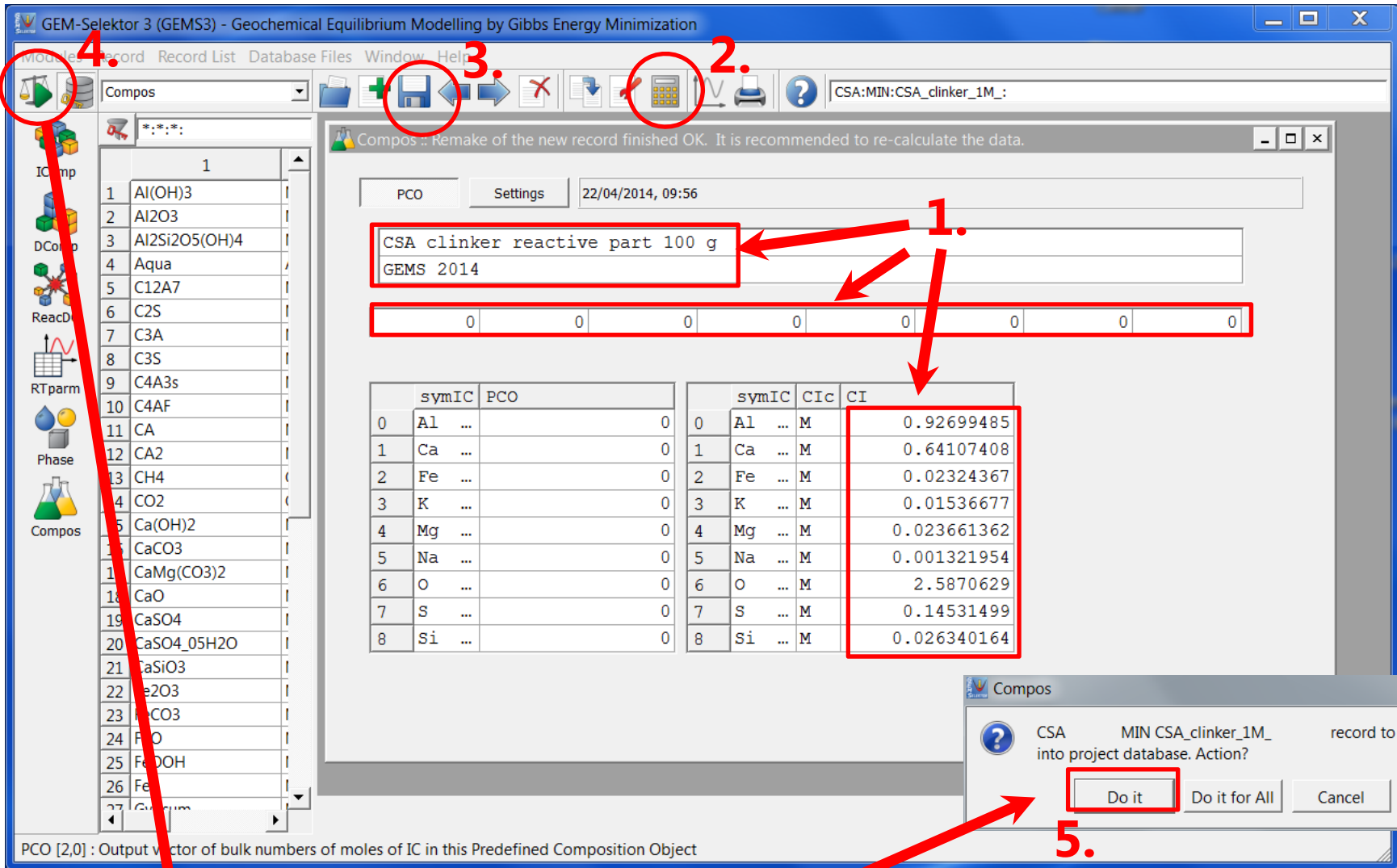
Learn more < Back **Finish** Cancel

5. Commercial CSA clinker + anhydrite + limestone



select elements

5. Commercial CSA clinker + anhydrite + limestone



1. CSA clinker reactive part 100 g
GEMS 2014

2. [Database Selection Icon]

3. [Save Icon]

4. [Load Icon]

5. Do it

	symIC	PCO		symIC	CIc	CI
0	Al	...	0	0	Al	M 0.92699485
1	Ca	...	0	1	Ca	M 0.64107408
2	Fe	...	0	2	Fe	M 0.02324367
3	K	...	0	3	K	M 0.01536677
4	Mg	...	0	4	Mg	M 0.023661362
5	Na	...	0	5	Na	M 0.001321954
6	O	...	0	6	O	M 2.5870629
7	S	...	0	7	S	M 0.14531499
8	Si	...	0	8	Si	M 0.026340164

Compos: Remake of the new record finished OK. It is recommended to re-calculate the data.

Compos: CSA MIN CSA_clinker_1M_ record to be inserted into project database. Action?

Do it Do it for All Cancel

PCO [2,0] : Output vector of bulk numbers of moles of IC in this Predefined Composition Object

load the project afterwards, your predefined composition will be inserted

5. Commercial CSA clinker + anhydrite + limestone



Open process file CSA_Cs_Cc_1

Input:

CSA:G:CSA_parent:0:0:1:20:0:CSA_Cs_Cc_1:S:

Controls Sampling Results Config 22/04/2014, 11:10

CSA clinker + anhydrite + calcite
GEMS 2014#

	iTm	iV	iP	iTC	iNv	iTau	ipXi	iNu	ipH	ipe
0	1000	0	1	20	0	0	0	0	0	0
1	1100	0	1	20	0	0	0	50	0	0
2	1	0	0	0	0	0	0	0.5	0	0
cTm	1100	0	1	20	0	0	0	50	0	0

```
$ hydration of 100 g solid with 100 g water and 1 g oxygen added
xa_{{Aqua}} =: 100;
xa_{{O2}} =: 1;
$
$ amount of calcium carbonate in g
xa_{{CaCO3}} =: 10;
$
$ amount of calcium sulfate is taken from iNu
xa_{{CaSO4}} =: cNu;
$
$ calculation of the amount of CSA
xa_{{CSA}} =: (100-xa_{{CaCO3}}-cNu)*0.756;
```

consideration of inert part

calculates with fixed CaCO_3 (here: 10 g) and varying ratio CSA/CaSO_4
(process file CSA_Cs_Cc_2 does it vice versa with fixed CaSO_4)

5. Commercial CSA clinker + anhydrite + limestone



Open process file CSA_Cs_Cc_1 Output (calculation of phase mass):

Controls	Sampling	Results	Config	22/04/2014, 11:10											
NeIt	9999	101	Next	0	I	0	J	100	Jp	100					
pSTkey				CSA:G:CSA_parent:0:0:1:20:0:				cTm	1100	cNV	0				
cTau		0		cpXi		0		cXi		1		cNu		50	
cpH		0		cpe		0		cEh		0		cT		293.15	

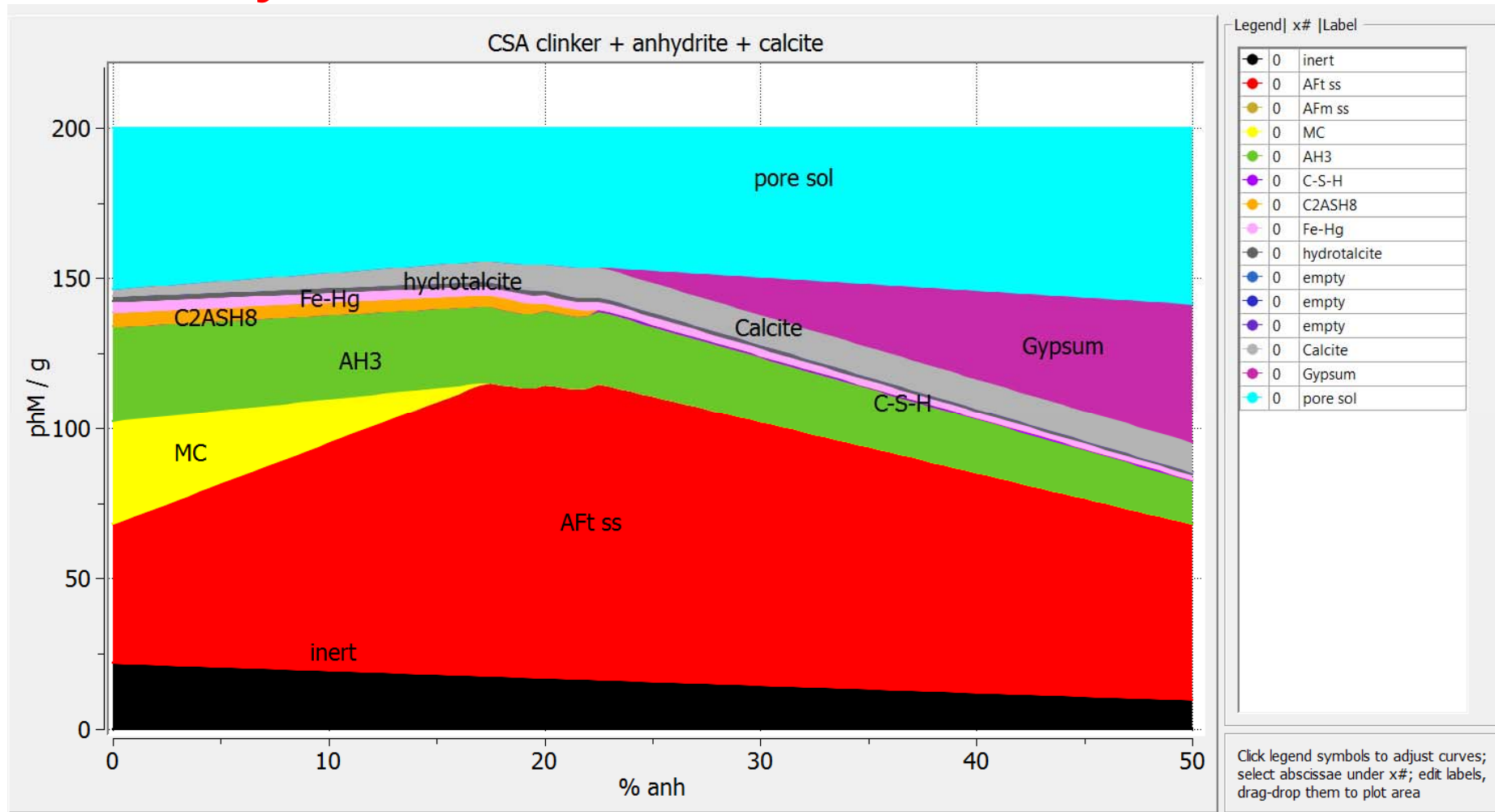

```
$ x-axis: amount of anhydrite added
xp[J] =: cNu;
$ v-axis: phase masses
yp[J][0] =: (100-xa [{CaCO3}]-cNu)*0.244;
yp[J][1] =: phM[{ettringite-Al}]+phM[{ettringite-Fe}]+phM[{SO4_CO3_Aft}]
+phM[{CO3_SO4_Aft}];
yp[J][2] =: phM[{monosulphate-Al}]+phM[{monosulphate-Fe}]
+phM[{SO4_OH_Afm}]+phM[{OH_SO4_Afm}];
yp[J][3] =: phM[{C4AcH11}];
yp[J][4] =: phM[{Al(OH)3mic}];
yp[J][5] =: phM[{CSHQ}];
yp[J][6] =: phM[{C2ASH8}];
yp[J][7] =: phM[{C3FS0.84H4.32}]+phM[{C3FS1.34H3.32}];
yp[J][8] =: phM[{OH-hydrotalcite}];
yp[J][9] =: 0;
yp[J][10] =: 0;
yp[J][11] =: 0;
yp[J][12] =: phM[{Calcite}];
yp[J][13] =: phM[{Gypsum}];
yp[J][14] =: phM[{aq_gen}];
```

consideration of inert part

5. Commercial CSA clinker + anhydrite + limestone



Result: ternary blend with 10% CaCO_3 and varying CSA/anhydrite



5. Commercial CSA clinker + anhydrite + limestone

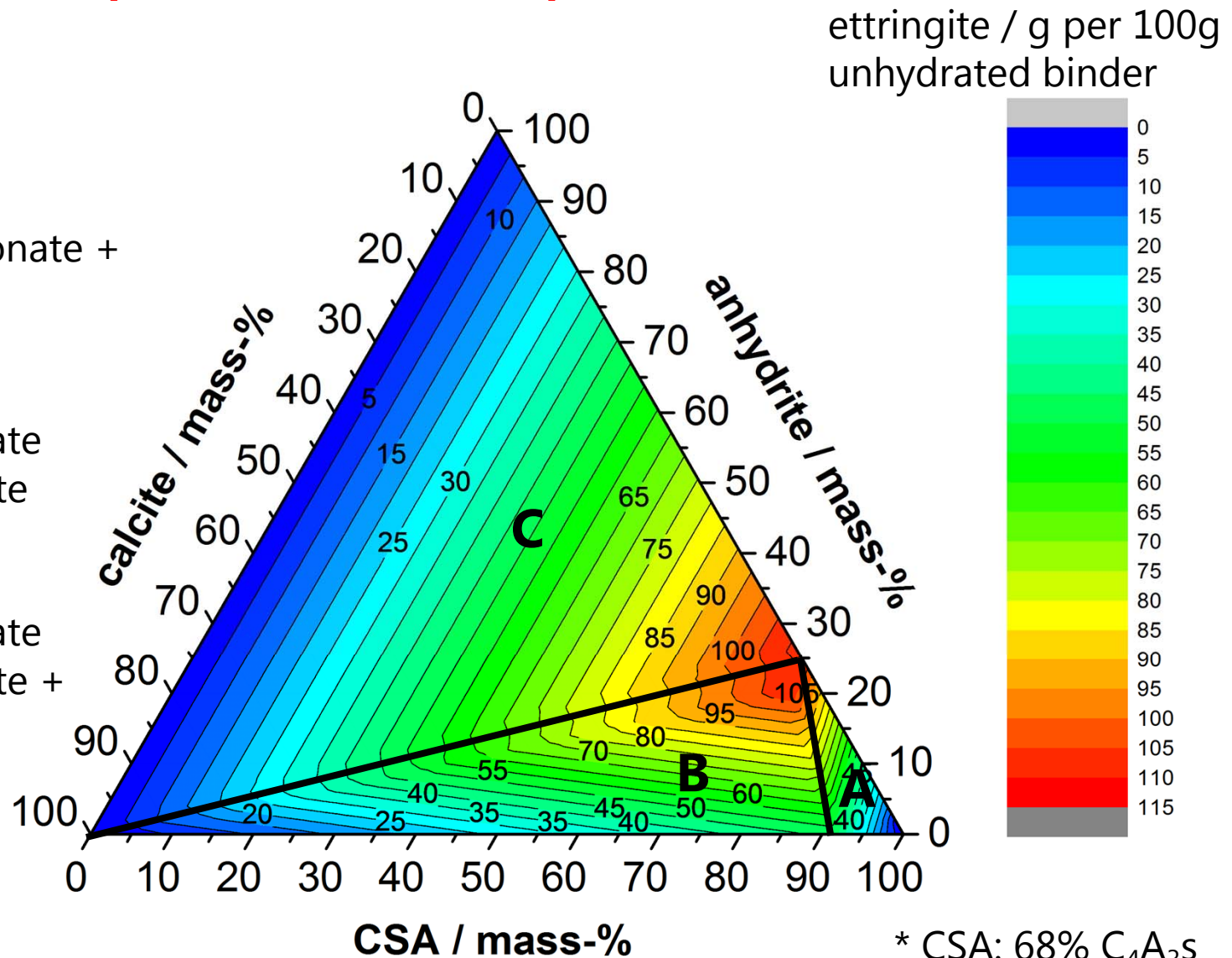


Phase diagram (commercial CSA*)

A: ettringite +
monosulfate
+ monocarbonate +
 AH_3

B: ettringite +
monocarbonate
+ AH_3 + calcite

C: ettringite +
monocarbonate
+ AH_3 + calcite +
gypsum



5. Commercial CSA clinker + anhydrite + limestone

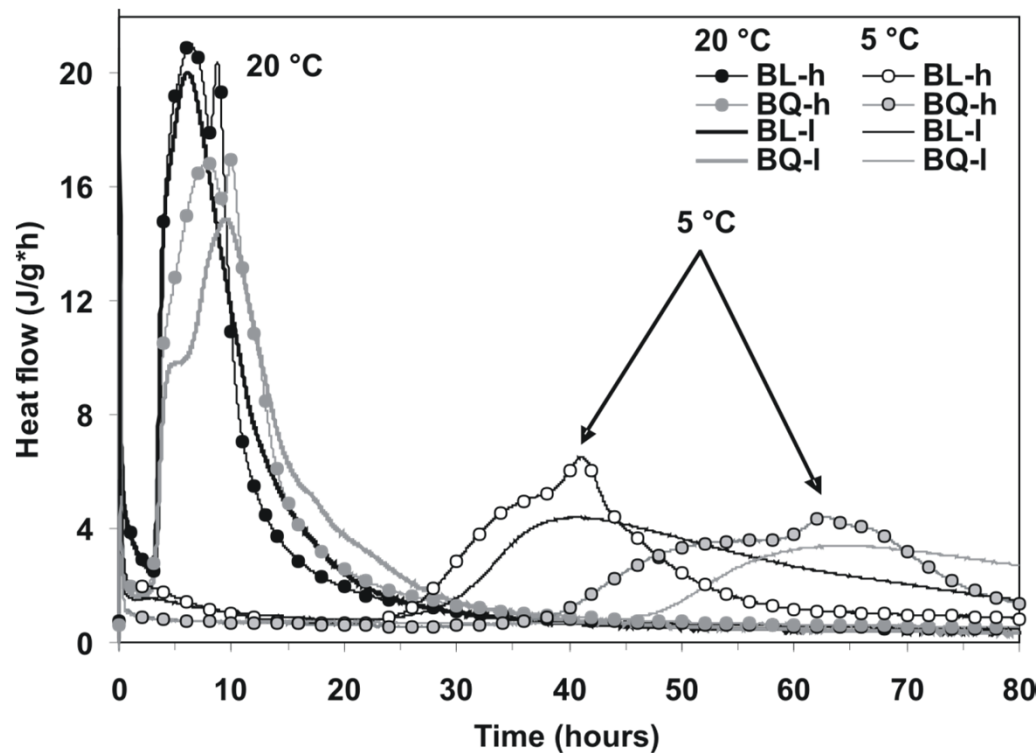


Conduction calorimetry

5 & 20°C, w/c = 0.80,
cement/filler \approx 1:1

Limestone powder

Quartz powder



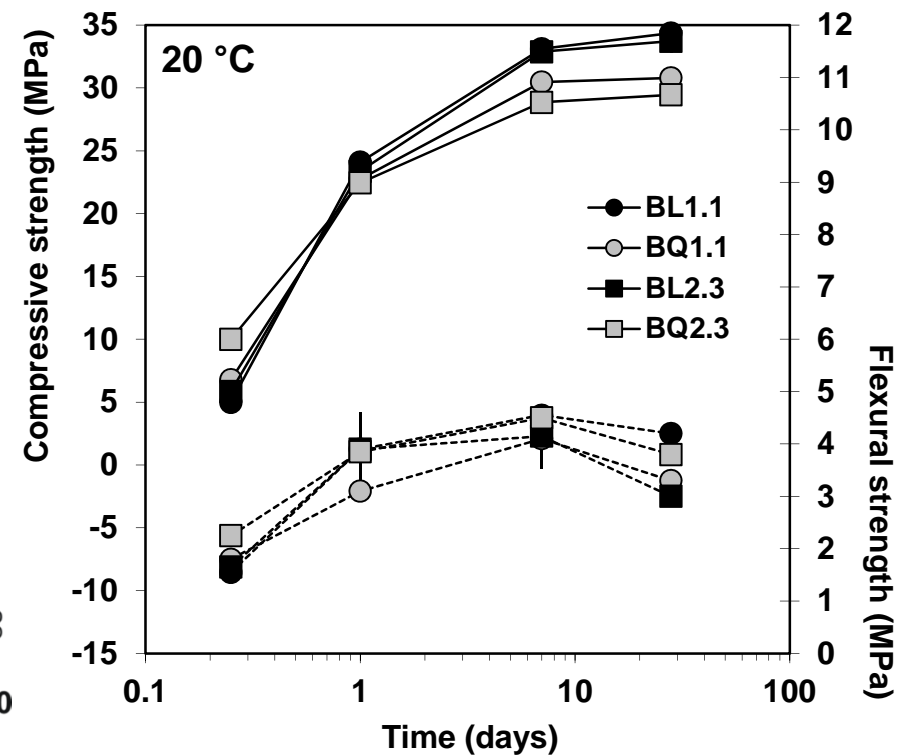
Compressive & flexural strength

20°C, w/c = 0.80

cement/filler/sand \approx 1:1:2

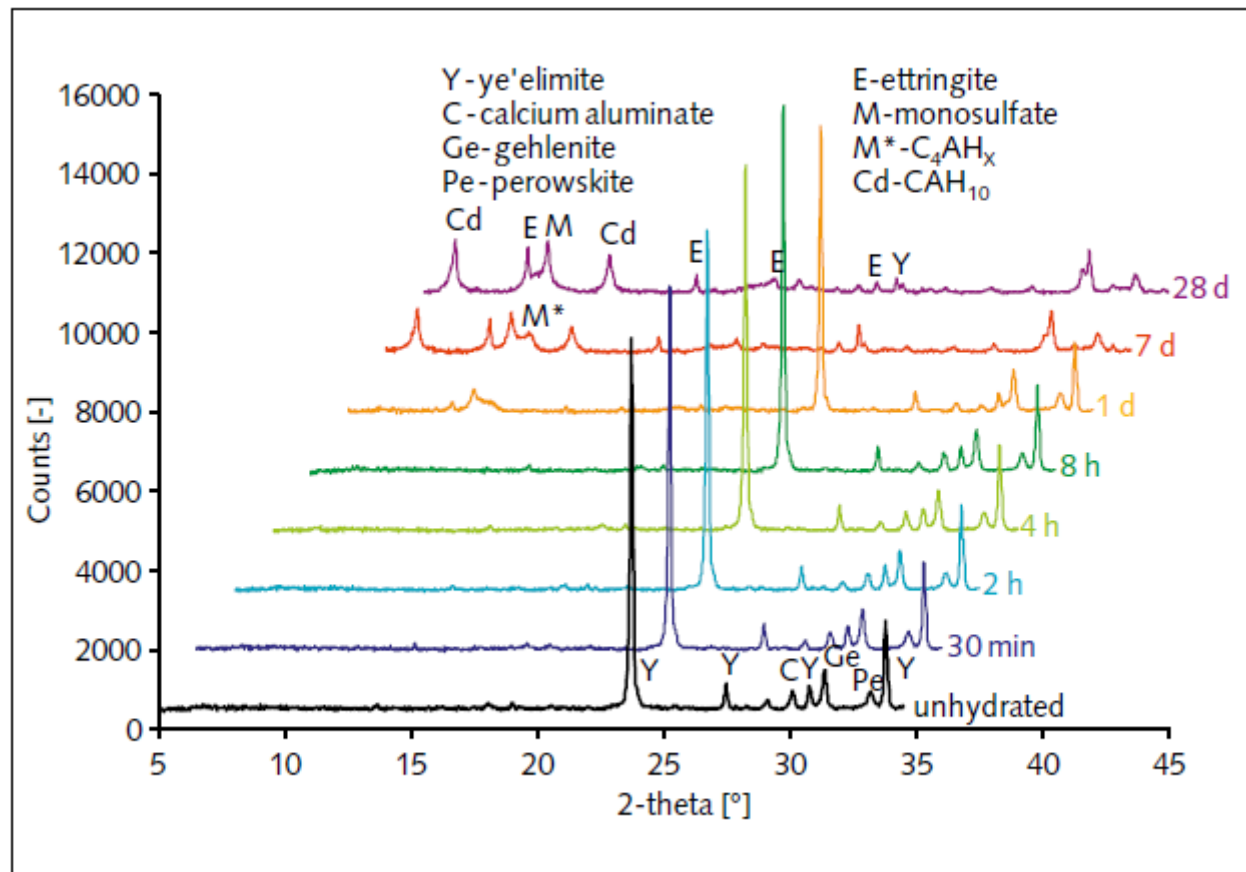
L = Limestone powder

Q = Quartz powder



6. Appendix: occurrence of CAH_{10}

CAH_{10} may occur as hydration product in CSA clinkers or CSA cements very low in calcium sulfate

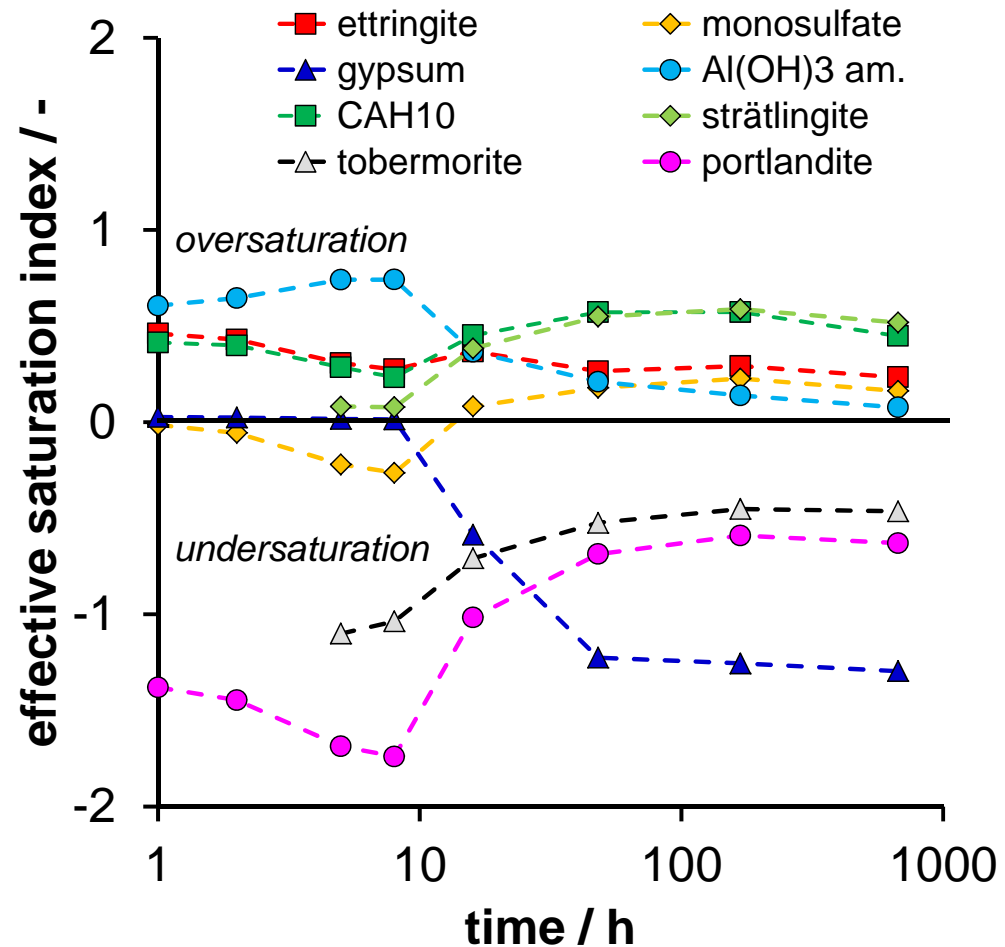


6. Appendix: occurrence of CAH_{10}

pore solution of CSA cements
may be oversaturated with
respect to CAH_{10}

The presence of CAH_{10} is linked
to the solubility of AH_3 .

AH_3 with high solubility:
=> CAH_{10} may form (as
predicted by GEMS)



6. Appendix: occurrence of CAH₁₀



project „Yeelimate“, single file „ye_anh“

Input: System Definition		Results: Equilibrium State										
Phase/species	L	T	On/	UC	Add to BC	UG	G0 corr.	UK	Lower_KC	Upper_KC	KC type	
aq_gen	29	a	+	g	0	J	0					
gas_gen	5	g	+	g	0	J	0					
SO4_OH_AFm	2	s	+	g	0	J	0					
OH_SO4_AFm	2	s	+	g	0	J	0					
SO4_CO3_Aft	2	s	-	g	0	J	0					
CO3_SO4_Aft	2	s	-	g	0	J	0					
Al(OH)3mic	1	s	+	g	0	J	0					
AlOHmic		O	+	M	0	J	5708	M	0	1e+006	B	
Gibbsite	1	s	-	g	0	J	0					
Graphite	1	s	-	g	0	J	0					
Mayenite	1	s	+	g	0	J	0					
Aluminate	1	s	+	g	0	J	0					
CA	1	s	+	g	0	J	0					
CA2	1	s	+	g	0	J	0					
C2AH75	1	s	+	g	0	J	0					
C3AH6	1	s	+	g	0	J	0					
C4AH13	1	s	-	g	0	J	0					
C4AH19	1	s	+	g	0	J	0					
CAH10	1	s	+	g	0	J	0					
C4AsH12	1	s	-	g	0	J	0					
C4Ac0.5H12	1	s	-	g	0	J	0					
C4AcH11	1	s	-	g	0	J	0					
ettringite	1	s	+	g	0	J	0					
Aragonite	1	s	-	g	0	J	0					
Calcite	1	s	-	g	0	J	0					
lime	1	s	+	g	0	J	0					
Portlandite	1	s	+	g	0	J	0					
Anhydrite	1	s	+	g	0	J	0					
Gypsum	1	s	+	g	0	J	0					
hemihydrate	1	s	+	g	0	J	0					

increases solubility product
by 1 log unit at 25°C

6. Appendix: occurrence of CAH₁₀



**Result: change of phase assemblage
CAH₁₀ instead of AFm**

Input: System Definition		Results: Equilibrium State	
Phase/species	L	T	Amount (mol)
+ a aq_gen	23	a	2.1481191
+ g gas_gen	3	g	0.03119591
+ s SO4_OH_AFm	2	s	0.10938991
+ s OH_SO4_AFm	2	s	2.8306647e-007
+ s Al(OH)3mic	1	s	0.59081956
+ s Mayenite	1	s	0
+ s Aluminate	1	s	0
+ s CA	1	s	0
+ s CA2	1	s	0
+ s C2AH75	1	s	0
+ s C3AH6	1	s	0
+ s C4AH19	1	s	0
+ s CAH10	1	s	0
+ s ettringite	1	s	0.037596495
+ s lime	1	s	0
+ s Portlandite	1	s	0
+ s Anhydrite	1	s	0
+ s Gypsum	1	s	0
+ s hemihydrate	1	s	0
+ s Sulphur	1	s	0



Input: System Definition		Results: Equilibrium State		
Phase/species		L	T	Amount (mol)
+	a aq_gen	23	a	0.53765046
+	g gas_gen	3	g	0.031237305
+	s SO4_OH_AFm	2	s	0
+	s OH_SO4_AFm	2	s	0
+	s Al(OH)3mic	1	s	0.29458899
+	s Mayenite	1	s	0
+	s Aluminate	1	s	0
+	s CA	1	s	0
+	s CA2	1	s	0
+	s C2AH75	1	s	0
+	s C3AH6	1	s	0
+	s C4AH19	1	s	0
+	s CAH10	1	s	0.22148015
+	s ettringite	1	s	0.073643289
+	s lime	1	s	0
+	s Portlandite	1	s	0
+	s Anhydrite	1	s	0
+	s Gypsum	1	s	0
+	s hemihydrate	1	s	0
+	s Sulphur	1	s	0

References



- Winnefeld F., Barlag S.: ZKG International 62 (2009), 42.
- Winnefeld F., Barlag S.: J. Therm. Anal. Calorim. 101 (2010), 949.
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- Pelletier-Chaignat L., Winnefeld F., Lothenbach B., Müller C.J.: Constr. Build. Mater. 26 (2012), 619.
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- Winnefeld F., Lothenbach B.: 1st International Conference on Sulphoaluminate Cement: Materials and Engineering Technology, Wuhan, China, October 23-24, 2013, 212.
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