

## Supplementary Information

### Effluent of Cooking with Active Oxygen and Solid Alkali (CAOSA): Components Separation, Recovery and Characterization

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## S1. Figures

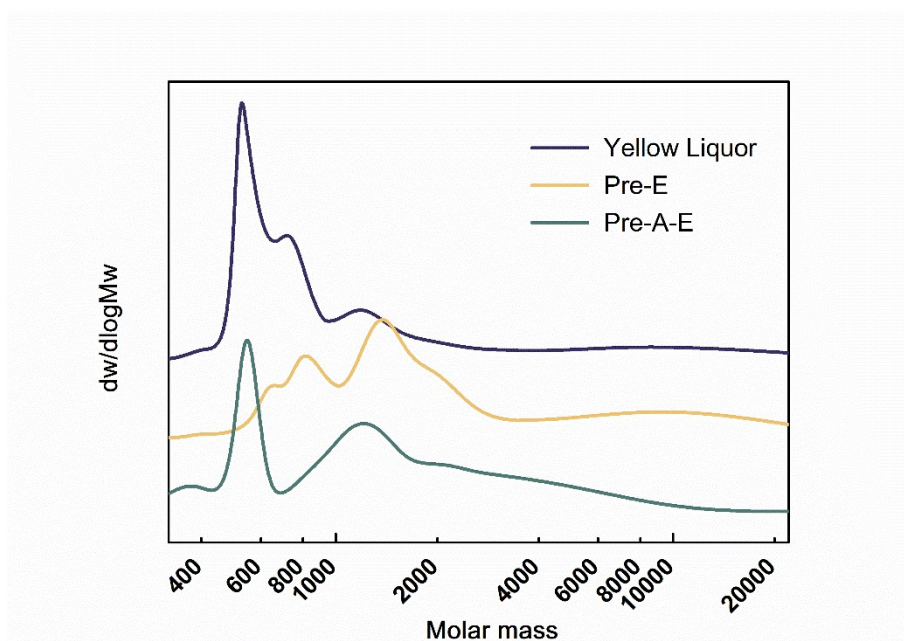


Figure S1 GPC results of YL/Pre-E/Pre-A-E

## S2. Tables

**Table S1 Dry components and precipitates in YL**

Entry	Dry components in YL/%	Precipitates in YL dry components/%			
		Pre-E	Pre-E-A	Pre-A	Pre-A-E
1	10.52	82.83	1.59	15.55	4.31
2	11.13	81.77	1.66	15.47	4.23
3	10.73	81.36	1.81	15.49	4.33
4	10.99	81.64	1.73	15.44	4.34
5	10.39	81.73	1.81	15.41	4.28
Average	10.75	81.87	1.72	15.47	4.30

**Table S2 Ash contents of YL dry components and precipitates**

Entry	Ash content/%				
	YL	Pre-E	Pre-E-A	Pre-A	Pre-A-E
1	25.23	27.03	3.35	1.63	0.81
2	25.77	27.58	3.78	1.64	0.88
3	25.06	28.21	3.82	1.7	0.84
Average	25.35	27.61	3.65	1.66	0.84

**Table S3 Assignments of main  $^{13}\text{C}$ - $^1\text{H}$  cross-signals in the HSQC spectra of****MWL**

Label	$\delta_{\text{C}}/\delta_{\text{H}}$ (ppm)	Assignments
$\text{C}_{\beta}$	53.1/3.46	$\text{C}_{\beta}$ - $\text{H}_{\beta}$ in phenylcoumaran (C)
$\text{B}_{\beta}$	53.5/3.07	$\text{C}_{\beta}$ - $\text{H}_{\beta}$ in $\beta$ - $\beta$ (resinol) (B)
$\text{OCH}_3$	56.4/3.70	C-H in methoxyls
$\text{A}_{\gamma}$	59.9/3.35-3.80	$\text{C}_{\gamma}$ - $\text{H}_{\gamma}$ in $\beta$ -O-4 substructures (A)
$\text{A}'_{\gamma}$	63.0/4.36	$\text{C}_{\gamma}$ - $\text{H}_{\gamma}$ in $\gamma$ -acylated $\beta$ -O-4 ( $\text{A}'$ )
$\text{C}_{\gamma}$	62.2/3.76	$\text{C}_{\gamma}$ - $\text{H}_{\gamma}$ in phenylcoumaran (C)
$\text{B}_{\gamma}$	71.2/3.82-4.18	$\text{C}_{\gamma}$ - $\text{H}_{\gamma}$ in $\beta$ - $\beta$ (resinol) (B)
$\text{A}_{\alpha}$	71.8/4.86	$\text{C}_{\alpha}$ - $\text{H}_{\alpha}$ in $\beta$ -O-4 unit (A)
$\text{A}_{\beta}(\text{G})$	83.4/4.38	$\text{C}_{\beta}$ - $\text{H}_{\beta}$ in $\beta$ -O-4 linked to G (A)
$\text{B}_{\alpha}$	84.8/4.66	$\text{C}_{\alpha}$ - $\text{H}_{\alpha}$ in $\beta$ - $\beta$ (resinol) (B)
$\text{A}'_{\beta}(\text{G})$	80.8/4.52	$\text{C}_{\beta}$ - $\text{H}_{\beta}$ in $\beta$ -O-4 linked to G ( $\text{A}'$ )
$\text{A}_{\beta}(\text{S})$	85.8/4.12	$\text{C}_{\beta}$ - $\text{H}_{\beta}$ in $\beta$ -O-4 linked to S (A)
$\text{C}_{\alpha}$	86.8/5.45	$\text{C}_{\alpha}$ - $\text{H}_{\alpha}$ in phenylcoumaran (C)
$\text{T}'_{2,6}$	103.9/7.34	$\text{C}'_{2,6}$ - $\text{H}'_{2,6}$ in triclin (T)
$\text{T}_6$	98.9/6.23	$\text{C}_{2,6}$ - $\text{H}_{2,6}$ in triclin (T)
$\text{T}_8$	94.2/6.60	$\text{C}_8$ - $\text{H}_8$ in triclin (T)
$\text{T}_3$	106.2/7.07	$\text{C}_3$ - $\text{H}_3$ in triclin (T)
$\text{S}_{2,6}$	103.9/6.70	$\text{C}_{2,6}$ - $\text{H}_{2,6}$ in syringyl unit (S)
$\text{S}'_{2,6}$	106.3/7.32	$\text{C}_{2,6}$ - $\text{H}_{2,6}$ in oxidized S unit (S)
$\text{G}_2$	110.8/6.97	$\text{C}_2$ - $\text{H}_2$ in guaiacyl unit (G)
$\text{G}_5$	114.5/6.70	$\text{C}_5$ - $\text{H}_5$ in guaiacyl unit (G)
$\text{G}_6$	119.0/6.78	$\text{C}_6$ - $\text{H}_6$ in guaiacyl unit (G)
$\text{H}_{2,6}$	127.7/7.17	$\text{C}_{2,6}$ - $\text{H}_{2,6}$ in p-hydroxyphenyl unit (H)
$\text{PCE}_{3,5}$	115.6/6.77	$\text{C}_{3,5}$ - $\text{H}_{3,5}$ in p-Coumarates (PCE)
$\text{PCE}_{2,6}$	130.2/7.48	$\text{C}_{2,6}$ - $\text{H}_{2,6}$ in p-Coumarates (PCE)

PCE <sub>7</sub>	144.8/7.51	C <sub>7</sub> -H <sub>7</sub> in p-Coumarates (PCE)
PCE <sub>8</sub>	113.7/6.24	C <sub>8</sub> -H <sub>8</sub> in p-Coumarates (PCE)
FA <sub>2</sub>	110.7/7.35	C <sub>2</sub> -H <sub>2</sub> in p-Ferulate (FA)
FA <sub>6</sub>	123.1/7.20	C <sub>6</sub> -H <sub>6</sub> in p-Ferulate (FA)
FA <sub>7</sub>	144.8/7.51	C <sub>7</sub> -H <sub>7</sub> in p-Ferulate (FA)

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**Table S4 Band assignment of lignin in FTIR spectra**

Entry	Wavenumber(cm <sup>-1</sup> )	Functional group
1	3412-3460	O-H stretch
2	3000-2842	C-H stretch in methyl and methylene groups
3	2900	C-H aliphatic axial deformation
4	2865	C-H of methoxy group
5	1738-1709	C=O stretch in unconjugated ketone, carbonyl and in ester groups; conjugated aldehydes and carboxylic acids absorb around and below 1700 cm-1
6	1655-1675	C=O stretch ; in conjugated p-subst. aryl ketones; strong electronegative substituents lower the wavenumber
7	1593-1605	aromatic skeletal vibrations plus C=O stretch; S > G; G condensed > G etherified
8	1505-1515	aromatic skeletal vibrations; G > S
9	1460-1470	C-H deformations; asym. in -CH3 and -CH2-
10	1422-1430	aromatic skeletal vibrations combined with C-H in-plane deform
11	1365-1370	aliphatic C-H stretch in CH3, not in OMe; phen. OH
12	1325-1330	S ring plus G ring condensed
13	1266-1270	G ring plus C=O stretch
14	1221-1230	C-C plus C-O plus C=O stretch; G condensed > G etherified
15	1166	typical for HGS lignins; C=O in ester groups (conj.)
16	1140	aromatic C-H in-plane deformation; typical for G units; whereby G condensed > etherified (typical

		for S units); plus secondary alcohol plus C=O stretch
17	1086	C-O deformation in secondary alcohols and aliphatics ethers
		aromatic C-H in-plane deformation, G > S; plus
18	1030-1035	C-O deform, in primary alcohols; plus C=O stretch (conj.)
19	966-990	-CH=CH- out-of-plane deform
20	915-925	C-H out-of-plane; aromatic
21	853-858	C-H out-of-plane in position 2,5, and 6 of G units
22	834-835	C-H out-of-plane in position 2,6 of S and in all position of H units
23	817-832	C-H out-of-plane in positions 2,5 and 6 of G units

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