

## Supporting Information

### Room-temperature conversion of CO<sub>2</sub> into quinazoline-2,4(1H,3H)-dione by deep eutectic solvents at atmospheric pressure with high efficiency

Yu Chen<sup>\*a</sup>, Chong Liu<sup>a</sup>, Yaoting Duan<sup>a</sup>, Dongkun Yu<sup>b</sup>, Zhenghui Liu<sup>c</sup>, Yuting Li<sup>a</sup>, Ruifen Shi<sup>b</sup>, Yuting Guo<sup>a</sup>, Tiancheng Mu<sup>\*b</sup>

<sup>a</sup> Department of Chemistry and Material Science, Langfang Normal University, Langfang 065000, Hebei, P.R. China. E-mail: yuchen@iccas.ac.cn; Phone: +86-316-2188211; Fax: +86-316-2112462. Zip: 065000.



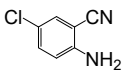
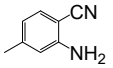
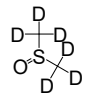
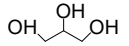

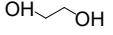
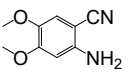
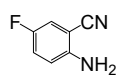
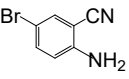
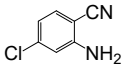
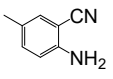
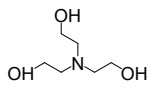
<sup>b</sup> Department of Chemistry, Renmin University of China, Beijing 100872, P.R. China. E-mail: tcmu@ruc.edu.cn; Phone: +86-10-62514925; Fax: +86-10-62516444. Zip: 100872.

<sup>c</sup> Department of Chemistry, Taizhou University, Taizhou 318000, Zhejiang, P. R. China. E-mail: liuzhenghui@126.com; Tel: +86-576-88661988; Fax: +86-576-88661988. Zip: 318000.

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Table S1. Structure, name, purity and CAS of chemicals investigated.

NO.	Structure	Name	Purity	CAS
1		2-aminobenzonitrile	98%	1885-29-6
2		1,5-diazabicyclo[4.3.0]non-5-ene (DBN)	98%	3001-72-7
3		2-amino-5-chlorobenzonitrile	98%	5922-60-1
4		2-amino-4-methylbenzonitrile	98%	26830-96-6
5		dimethylsulfoxide-d6 (d-DMSO)	99.8 atom % D	2206-27-1
6		glycerol (GLY)	99%	56-81-5
7		methyl tert-butyl ether	99%	1634-04-4
8		ethylene glycol (EG)	99%	107-21-1
9		2-amino-4,5-dimethoxybenzonitrile	98%	26961-27-3
10		2-amino-5-fluorobenzonitrile	98%	61272-77-3
11		2-amino-5-bromobenzonitrile	98%	39263-32-6
12		2-amino-4-chlorobenzonitrile	98%	38487-86-4
13		2-amino-5-methylbenzonitrile	98%	5925-93-9
14		triethanolamine (TEA)	99%	102-71-6

**Table S2. Fitted parameters for the relationship between yields and time at 30 °C and 60 °C.<sup>a</sup>**

<i>T</i> / °C	<i>a</i> / %	<i>b</i> / h <sup>-1</sup>	<i>R</i> <sup>2</sup>
30	50.2	0.0618	0.9987
60	100	0.1189	0.9234

<sup>a</sup>  $y = a(1 - e^{-bt})$ , *y* and *t* represent yield and time, respectively. *a* and *b* are the fitted parameters.