

# Electronic Supplementary Materials

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## Mass transfer enhancement of hollow fiber membrane deoxygenation by Dean vortices

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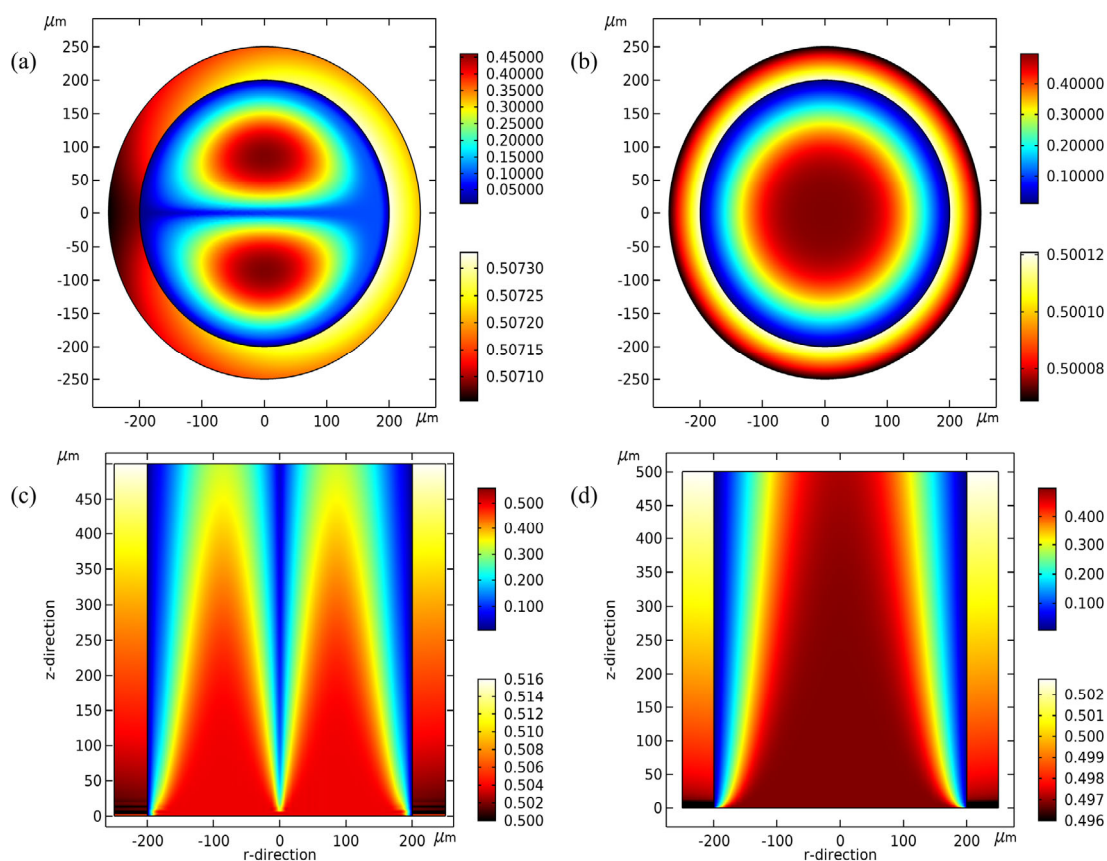
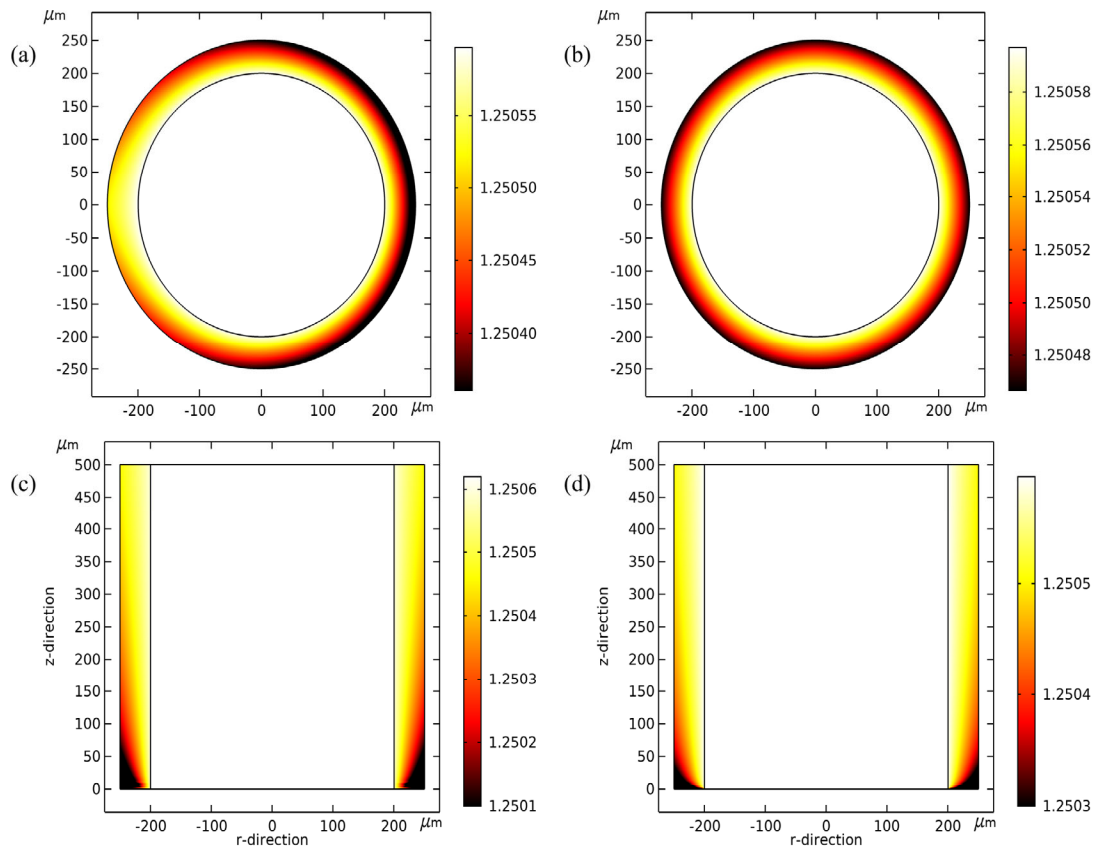


Fig. S1 Concentration profiles of N<sub>2</sub> of (a) HFFM deoxygenation ( $\kappa=162.9 \text{ m}^{-1}$ ,  $z=0.5 \text{ m}$ ), (b) SHFM deoxygenation ( $\kappa=0$ ,  $z=0.5 \text{ m}$ ) on the cross-section and (c) HFFM deoxygenation ( $\kappa=162.9 \text{ m}^{-1}$ ), (d) SHFM deoxygenation ( $\kappa=0$ ) along the axis direction.  $Re=217$ ,  $T=298.2 \text{ K}$ ,  $\bar{v}=0.5 \text{ m s}^{-1}$ ,  $p/p_0=0.50$ ,  $l=1.0 \text{ m}$



**Fig. S2** Porous membrane concentration profiles of  $\text{H}_2\text{O}$  vapor of (a) HHF membrane deoxygenation ( $\kappa = 162.9 \text{ m}^{-1}$ ,  $z = 0.5 \text{ m}$ ), (b) SHF membrane deoxygenation ( $\kappa = 0$ ,  $z = 0.5 \text{ m}$ ) on the cross-section and (c) HHF membrane deoxygenation ( $\kappa = 162.9 \text{ m}^{-1}$ ), (d) SHF membrane deoxygenation ( $\kappa = 0$ ) along the axis direction.  $Re = 217$ ,  $T = 298.2 \text{ K}$ ,  $\bar{v} = 0.5 \text{ m s}^{-1}$ ,  $p/p_0 = 0.50$ ,  $l = 1.0 \text{ m}$

**Table S1 Parameters used for simulation**

$T$ (K)	$D_{O_2,l}$ ( $\times 10^{-9} \text{ m}^2 \text{ s}^{-1}$ )	$D_{N_2,l}$ ( $\times 10^{-9} \text{ m}^2 \text{ s}^{-1}$ )	$D_{O_2,k}$ ( $\times 10^{-5} \text{ m}^2 \text{ s}^{-1}$ )	$D_{N_2,k}$ ( $\times 10^{-5} \text{ m}^2 \text{ s}^{-1}$ )	$D_{H_2O,k}$ ( $\times 10^{-5} \text{ m}^2 \text{ s}^{-1}$ )	$\mu_w$ ( $\times 10^{-6} \text{ m}^2 \text{ s}^{-1}$ )
278.2	2.011	1.254	2.859	3.057	3.812	1.377
288.2	2.084	1.495	2.910	3.111	3.880	1.129
298.2	2.156	1.803	2.960	3.165	3.946	0.919
308.2	2.228	2.176	3.010	3.217	4.012	0.748

**Table S1 (continued) Parameters used for simulation**

$T$ (K)	$\mu_{O_2,g}$ ( $\times 10^{-5} \text{ Pa s}$ )	$\mu_{N_2,g}$ ( $\times 10^{-5} \text{ Pa s}$ )	$\mu_{H_2O,g}$ ( $\times 10^{-6} \text{ Pa s}$ )	$H_{O_2}$ (GPa)	$H_{N_2}$ (GPa)	$p_w$ (kPa)
278.2	1.945	1.682	8.214	2.893	6.001	0.847
288.2	1.998	1.726	8.614	3.692	7.446	1.685
298.2	2.050	1.769	9.014	4.423	8.752	3.098
308.2	2.103	1.813	9.414	5.086	9.918	5.555

**Table S2** Experimental data of HHFM and SHFM deoxygenation

No.	$\nu$ (m s <sup>-1</sup> )	$\kappa$ (m <sup>-1</sup> )	$p/p_0$	$T$ (K)	$c_f$ (mol m <sup>-3</sup> )	$c_{out}$ (mol m <sup>-3</sup> )	$\sigma$ (x10 <sup>-2</sup> )*
1	0.10	162.9	0.90	298.2	0.256	0.246	1.81
2	0.25	162.9	0.90	298.2	0.256	0.249	1.33
3	0.50	162.9	0.90	298.2	0.256	0.252	2.43
4	0.75	162.9	0.90	298.2	0.256	0.253	1.11
5	1.00	162.9	0.90	298.2	0.256	0.253	1.93
6	1.25	162.9	0.90	298.2	0.256	0.254	2.69
7	0.10	162.9	0.70	298.2	0.256	0.224	0.59
8	0.25	162.9	0.70	298.2	0.256	0.236	2.18
9	0.50	162.9	0.70	298.2	0.256	0.244	2.52
10	0.75	162.9	0.70	298.2	0.256	0.247	1.98
11	1.00	162.9	0.70	298.2	0.256	0.249	2.07
12	1.25	162.9	0.70	298.2	0.256	0.251	1.53
13	0.10	162.9	0.50	298.2	0.256	0.188	0.55
14	0.25	162.9	0.50	298.2	0.256	0.212	2.74
15	0.50	162.9	0.50	298.2	0.256	0.227	2.35
16	0.75	162.9	0.50	298.2	0.256	0.233	1.36
17	1.00	162.9	0.50	298.2	0.256	0.236	1.28
18	1.25	162.9	0.50	298.2	0.256	0.239	1.95
19	0.10	162.9	0.30	298.2	0.256	0.139	1.37
20	0.25	162.9	0.30	298.2	0.256	0.178	1.51
21	0.50	162.9	0.30	298.2	0.256	0.203	1.61
22	0.75	162.9	0.30	298.2	0.256	0.213	2.08
23	1.00	162.9	0.30	298.2	0.256	0.219	2.57
24	1.25	162.9	0.30	298.2	0.256	0.222	1.83
25	0.10	162.9	0.050	298.2	0.256	0.0550	0.36
26	0.25	162.9	0.050	298.2	0.256	0.118	1.46

27	0.50	162.9	0.050	298.2	0.256	0.159	1.44
28	0.75	162.9	0.050	298.2	0.256	0.176	0.74
29	1.00	162.9	0.050	298.2	0.256	0.185	2.49
30	1.25	162.9	0.050	298.2	0.256	0.192	2.18
31	0.10	0	0.90	298.2	0.256	0.248	2.80
32	0.25	0	0.90	298.2	0.256	0.252	1.43
33	0.50	0	0.90	298.2	0.256	0.254	1.58
34	0.75	0	0.90	298.2	0.256	0.255	2.03
35	1.00	0	0.90	298.2	0.256	0.255	0.88
36	1.25	0	0.90	298.2	0.256	0.255	1.91
37	0.10	0	0.70	298.2	0.256	0.232	2.39
38	0.25	0	0.70	298.2	0.256	0.243	1.83
39	0.50	0	0.70	298.2	0.256	0.250	1.98
40	0.75	0	0.70	298.2	0.256	0.252	1.36
41	1.00	0	0.70	298.2	0.256	0.253	2.53
42	1.25	0	0.70	298.2	0.256	0.254	1.47
43	0.10	0	0.50	298.2	0.256	0.205	1.46
44	0.25	0	0.50	298.2	0.256	0.227	1.78
45	0.50	0	0.50	298.2	0.256	0.240	1.74
46	0.75	0	0.50	298.2	0.256	0.245	2.18
47	1.00	0	0.50	298.2	0.256	0.248	2.15
48	1.25	0	0.50	298.2	0.256	0.249	2.53
49	0.10	0	0.30	298.2	0.256	0.169	1.61
50	0.25	0	0.30	298.2	0.256	0.205	1.57
51	0.50	0	0.30	298.2	0.256	0.226	2.31
52	0.75	0	0.30	298.2	0.256	0.235	1.90
53	1.00	0	0.30	298.2	0.256	0.240	1.06
54	1.25	0	0.30	298.2	0.256	0.243	2.40

55	0.10	0	0.050	298.2	0.256	0.104	0.10
56	0.25	0	0.050	298.2	0.256	0.166	0.47
57	0.50	0	0.050	298.2	0.256	0.201	2.11
58	0.75	0	0.050	298.2	0.256	0.217	1.83
59	1.00	0	0.050	298.2	0.256	0.225	1.94
60	1.25	0	0.050	298.2	0.256	0.230	2.46
61	0.10	162.9	0.050	278.2	0.391	0.111	0.83
62	0.25	162.9	0.050	278.2	0.391	0.203	1.32
63	0.50	162.9	0.050	278.2	0.391	0.259	1.62
64	0.75	162.9	0.050	278.2	0.391	0.282	1.22
65	1.00	162.9	0.050	278.2	0.391	0.297	0.49
66	1.25	162.9	0.050	278.2	0.391	0.309	0.59
67	0.10	162.9	0.050	288.2	0.314	0.0791	0.43
68	0.25	162.9	0.050	288.2	0.314	0.154	0.92
69	0.50	162.9	0.050	288.2	0.314	0.200	1.47
70	0.75	162.9	0.050	288.2	0.314	0.220	0.67
71	1.00	162.9	0.050	288.2	0.314	0.233	2.07
72	1.25	162.9	0.050	288.2	0.314	0.243	2.24
73	0.10	162.9	0.050	308.2	0.218	0.0352	0.24
74	0.25	162.9	0.050	308.2	0.218	0.0898	0.87
75	0.50	162.9	0.050	308.2	0.218	0.124	0.89
76	0.75	162.9	0.050	308.2	0.218	0.140	1.22
77	1.00	162.9	0.050	308.2	0.218	0.151	1.24
78	1.25	162.9	0.050	308.2	0.218	0.159	1.55
79	0.10	0	0.050	278.2	0.391	0.178	0.65
80	0.25	0	0.050	278.2	0.391	0.270	0.71
81	0.50	0	0.050	278.2	0.391	0.318	1.89
82	0.75	0	0.050	278.2	0.391	0.338	2.00

83	1.00	0	0.050	278.2	0.391	0.349	2.92
84	1.25	0	0.050	278.2	0.391	0.357	2.06
85	0.10	0	0.050	288.2	0.314	0.137	0.96
86	0.25	0	0.050	288.2	0.314	0.212	1.16
87	0.50	0	0.050	288.2	0.314	0.252	1.56
88	0.75	0	0.050	288.2	0.314	0.269	2.07
89	1.00	0	0.050	288.2	0.314	0.278	1.58
90	1.25	0	0.050	288.2	0.314	0.285	1.31
91	0.10	0	0.050	308.2	0.218	0.0842	0.52
92	0.25	0	0.050	308.2	0.218	0.139	0.92
93	0.50	0	0.050	308.2	0.218	0.169	2.03
94	0.75	0	0.050	308.2	0.218	0.182	0.68
95	1.00	0	0.050	308.2	0.218	0.189	1.97
96	1.25	0	0.050	308.2	0.218	0.194	1.51

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\*  $\sigma$  is the standard deviation of  $C_{out}$  obtained from three parallel experimental data.

**Table S3 Experimental results of long-time run for 100 h at  $Re=43.5$ ,  $T=298.2$  K,  $k=162.9$  m<sup>-1</sup>,  
 $\bar{v}=0.1$  ms<sup>-1</sup>,  $p/p_0=0.50$**

$t$ (h)	$c_{\text{out}}$ (mol m <sup>-3</sup> )
0.20	0.0550
0.50	0.0553
1.0	0.0550
5.0	0.0548
10	0.0550
25	0.0553
50	0.0553
75	0.0550
100	0.0548

**Table S4 Boundary conditions used to model Liu's (Liu *et al.*, 2005) and Kaufhold's (Kaufhold *et al.*, 2012) works.**

Mathematical expressions	
<i>Boundary conditions of Liu's work</i>	
Lumen side aqueous phase ( $i=O_2, N_2$ )	
$z = 0 : c_{O_2} = c_{O_2,f} \quad c_{N_2} = 0,$	(S1)
$z = l : \partial c_i / \partial z = 0,$	(S2)
$r = 0 : \partial c_i / \partial r = 0.$	(S3)
Lumen side-membrane pores interface ( $i=O_2, N_2$ )	
$r = r_1 : (D_i \partial c_i / \partial r) _{\text{lumen}} = N_i _{\text{membrane}} = (-\bar{D}_i \nabla c_i + c_i \bar{\mathbf{v}}) _{\text{membrane}},$	(S4)
$r = r_1 : (H_i c_i / (c_i + \rho_w / M_w) / RT) _{\text{lumen}} = c_i _{\text{membrane}},$	(S5)
$r = r_1 : c_{H_2O} _{\text{membrane}} = p_w.$	(S6)
Membrane pores-shell side interface ( $i=O_2, N_2, H_2O$ )	
$r = r_2 : c_{N_2} = c _{\text{shell}} \quad c_{O_2} = c_{H_2O} = 0.$	(S7)
<i>Boundary conditions of Kaufhold's work</i>	
Lumen side aqueous phase ( $i=O_2, N_2$ )	
$z = 0 : c_{N_2} = c_{N_2,f} \quad c_{O_2} = 0,$	(S8)
$z = l : \partial c_i / \partial z = 0,$	(S9)
$r = 0 : \partial c_i / \partial r = 0.$	(S10)
Lumen side-membrane pores interface ( $i=O_2, N_2$ )	
$r = r_1 : (D_i \partial c_i / \partial r) _{\text{lumen}} = N_i _{\text{membrane}} = (-\bar{D}_i \nabla c_i + c_i \bar{\mathbf{v}}) _{\text{membrane}},$	(S11)
$r = r_1 : (H_i c_i / (c_i + \rho_w / M_w) / RT) _{\text{lumen}} = c_i _{\text{membrane}},$	(S12)
$r = r_1 : c_{H_2O} _{\text{membrane}} = p_w.$	(S13)
Membrane pores-shell side interface ( $i=O_2, N_2, H_2O$ )	
$r = r_2 : (RT \sum c_i) _{\text{membrane}} = p _{\text{shell}},$	(S14)

$$r = r_2 : c_i / \sum c_i = N_i / \sum N_i. \quad (\text{S15})$$

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