



1	Article
1	лисие

2	Reductive dechlorination of chloroacetamides with
3	NaBH4 catalyzed by zero valent iron, ZVI, nanoparticles
4	in ORMOSIL matrices prepared via the sol-gel route
5	Michael Meistelman ¹ , Dan Meyerstein ^{2,3} *, Amos Bardea ⁴ , Ariela Burg ⁵ , Dror Shamir ^{6,} and Yael Albo ¹ *
6	¹ Chemical Engineering Dept. and The Centre for Radical Reactions, Ariel University, Ariel, Israel.
7	² Department of Chemical Sciences and The Centre for Radical Reactions, Ariel University, Ariel, Israel.
8	³ Chemistry Dept., Ben-Gurion University, Beer-Sheva, Israel
9	⁴ Department of Electrical and Electronics Engineering, Faculty of Engineering, Holon Institute of Technology (HIT),
10	Holon, Israel
11	⁵ Department of Chemical Engineering, Sami Shamoon College of Engineering, Beer-Sheva, Israel.
12 13 14 15	⁶ Department of Chemistry, Nuclear Research Centre Negev, Beer-Sheva, Israel. E-mail addresses: michaelme@ariel.ac.il (M. Meistelman), danm@ariel.ac.il (D. Meyerstein), amos.bardea@hit.ac.il, arielab@ac.sce.ac.il (A. Burg), drorshamir@gmail.com (D. Shamir) yaelyt@ariel.ac.il (Y. Albo)
16 17	Received: date; Accepted: date; Published: date
18	
19 20	
21	
22	
23	
24	
25	
26 27	
21 28	
29	
30	

5. Supplementary Materials:



Figure S1. (a) Raw wet ORMOSIL gel, (b) Crushed ZVI@ORMOSIL gel

- Table S 1 is a summary of survey scan peaks and their integrated areas following the surface

xps analysis of 1% ZVI@ORMOSIL

Table S 1: XPS surface analysis elemental composition of 1% ZVI@ORMOSIL

Name	Peak (BE)	FWHM	Area	At%
Fe 2p	714.40	4.53	53140	0.2
O 1s	535.16	3.30	4168746	57.6
C 1s	287.34	3.56	336692	11.2
Si 2p	106.02	3.34	933002	31.0

Formula Fe			d	20	l fix	h	k	I
Name Iron			2.02600	44.693	999	1	1	0
Name (mineral) Iron, syn			1.43260	65.053	126	2	0	0
Name (common) ±-Fe			1.16971	82.377	206	2	1	1
			1.01300	99.001	58	2	2	0
			0.90606	116.460	86	3	1	0
			0.82711	137.280	25	2	2	2
Lattice: Cubic	Mol. weight =	55.85						
S.G.: Im-3m (229)	Volume [CD] =	23.52						
	Dx =	7.88						
	Dm =							
a = 2.86520	l/lcor =	11.580						
a/h 7 = 0								
a/b 1.00000 Z = 2								
c/b 1 00000								
= 1.00000								
LPF Collection Code: 1503158 Sample Propagation: STARTING MAT		la envetale						
grown by chemical transport method.	transport agent iod	ie crystais ine						
Temperature of Data Collection: 297	ĸ							
Unit Cell Data Source: Single Crystal								
Radiation: CuK±1 Fill	ter: Not spe	cified						
Wavelength 1.54060 d-s	spacing:							
SS/EOM: 000.0 (0.0001.6)								
GGH GML 333.3 (0.0001,0)								

Pattern: PDF 04-007-9753 Radiation: 1.54060 Quality: Star (*)

48 Figure S2. Fe⁰ phases powder diffraction file

- 4

Formula Fe3 O4		d	20	l fix	h	k	1
Name Iron Oxide		4.84732	18.288	84	1	1	1
Name (mineral) Magnetite,	syn	2.96836	30.081	286	2	2	0
Name (common) iron diiron(III) oxide	2.53143	35.431	1000	3	1	1
		2.42366	37.063	78	2	2	2
		2.09895	43.061	206	4	0	0
		1.92613	47.146	6	3	3	1
Lattice: Cubic	Mol. weight = 231.53	1.71379	53.419	88	4	2	2
S.G.: Fd-3m (227)	Volume [CD] = 591.82	1.61577	56.945	282	5	1	1
	Dx = 5.2	1.48418	62.531	377	4	4	0
		1.41915	65.747	8	5	3	1
a = 8.39580	5.100	1.39930	66.801	1	4	4	2
		1.32749	70.939	29	6	2	0
a/b		1.28035	73.974	72	5	3	3
= 1.00000		1.26571	74.975	31	6	2	2
c/b 1.00000		1.21183	78.936	22	4	4	4
-		1.17565	81.871	4	7	1	1
		1.12194	86.720	30	6	4	2
ANY: A2Y4		1.09304	89.616	105	7	3	1
LPF Collection Code: 541353		1.04947	94.444	39	8	0	0
Sample Preparation: STARTING MA	TERIALS:Fe2O3,Fe	1.02571	97.353	1	7	3	3
silica tube. Sample annealed at 107	3 K for 95 d	1.01814	98.326	1	6	4	4
Unit Cell Data Source: Powder Diffra	action	0.98945	102.248	14	8	2	2
		0.96946	105.228	57	7	5	1
		0.96306	106.230	12	6	6	2
		0.93868	110.294	23	8	4	0
		0.92156	113.412	2	9	1	1
		0.91606	114.467	1	8	4	2
		0.89499	118.785	7	6	6	4
		0.88012	122.142	42	9	3	1
		0.85689	128.039	89	8	4	4
		0.84381	131.813	1	9	3	3
		0.82328	138.667	19	10	2	0
		0.81165	143.263	56	9	5	1
		0.80789	144.908	13	10	2	2
Radiation: CuK±1 F Wavelength 1.54060 d- : SS/FOM: 999.9 (0.0001,30)	ilter: Not specified spacing:						

Pattern: PDF 04-005-4319 Radiation: 1.54060 Quality: Star (*)

57 Figure S3. Fe₃O₄ phases powder diffraction file

58



(a)

(b)

Figure S4.Y-axis dV/dD (cm³g⁻¹nm⁻¹), X-axis Diameter (nm) (a) pore size distribution for Blank@Ormosil, (b) pore size
distribution for 1% ZVI@ORMOSIL



© 2020 by the authors. Submitted for possible open access publication under the terms and conditions of the Creative Commons Attribution (CC BY) license (http://creativecommons.org/licenses/by/4.0/).