

# Facile Green Synthesis of Wasted Hops-Based Zinc Oxide Nanozymes as Peroxidase-Like Catalysts for Colorimetric Analysis

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## 1. Additional Figures

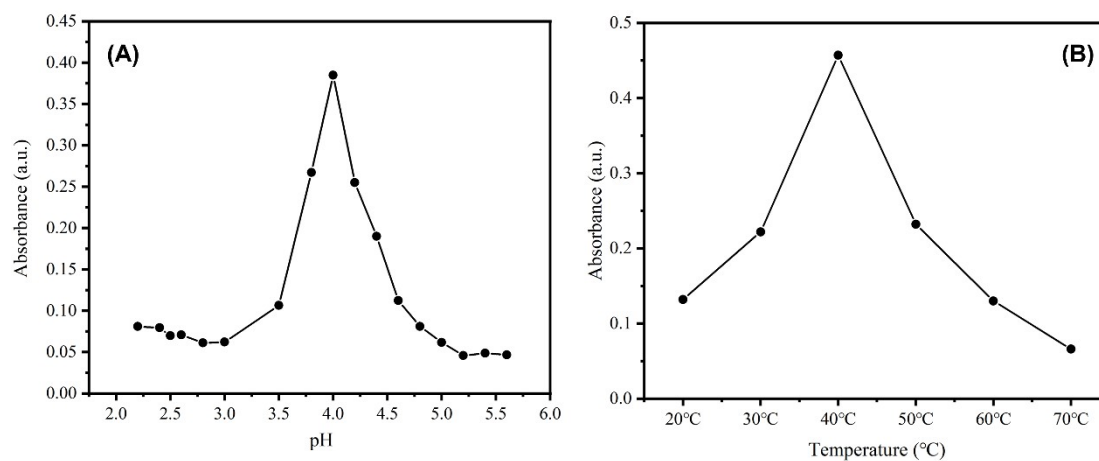


Figure S1. (A) pH-dependent, (B) temperature-dependent POD-like activity of WHE-ZnO NEs.

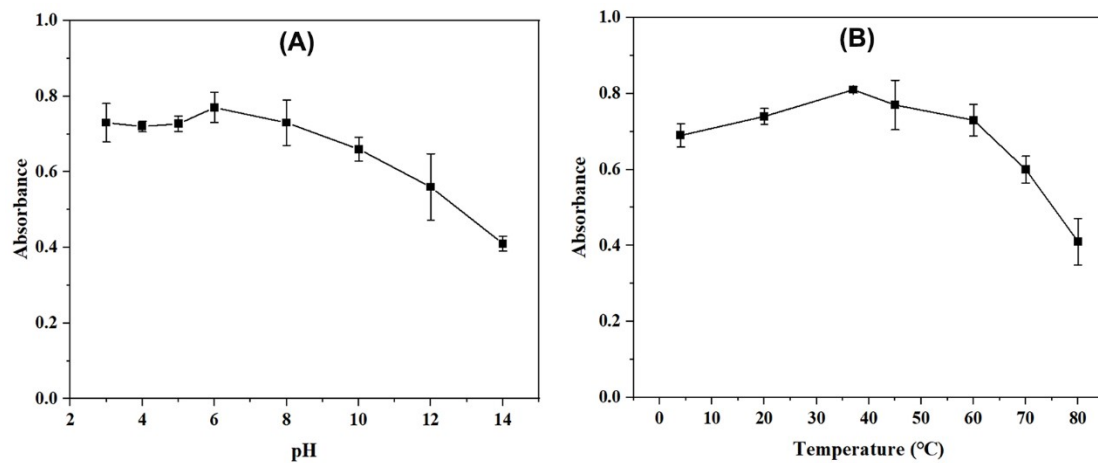


Figure S2. Stability of WHE-ZnO NEs after incubating at various temperature (A) and pH values (B) for 24 h.

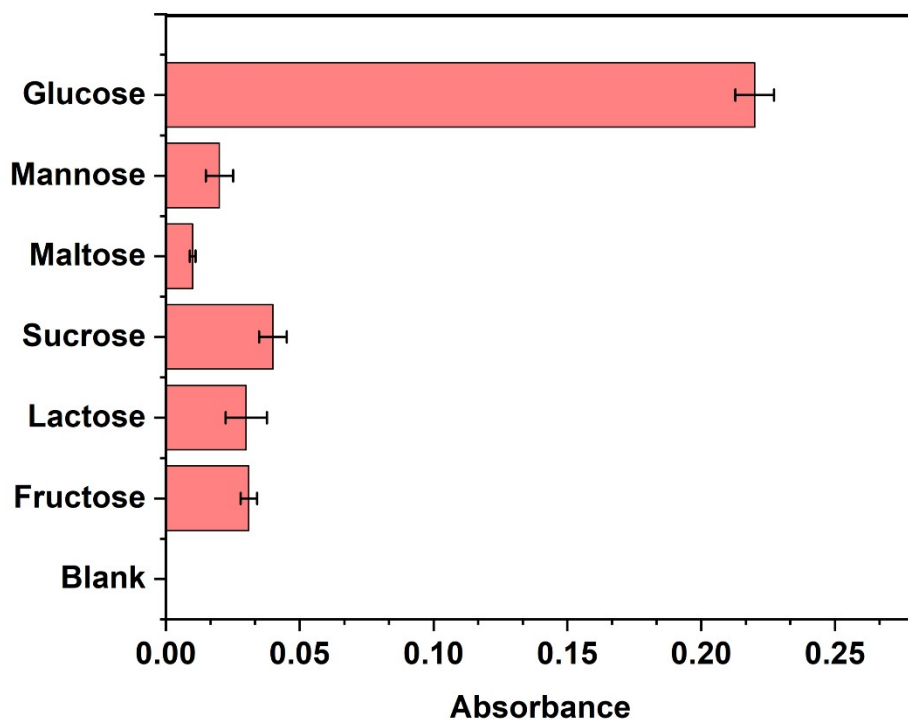


Figure S3. Selectivity evaluation of the glucose detection method.

## 2. Additional Tables

Table S1. Analysis of active ingredients in hops extract

No.	RT (min)	Compound	Molecular formula	Relative content (%)
1	5.42	Ribitol, 1,3:2,4-di-O-benzylidene	C <sub>19</sub> H <sub>20</sub> O <sub>5</sub>	0.13
2	6.10	1,5-Heptadiene, 3,3,6-trimethyl	C <sub>10</sub> H <sub>18</sub>	4.04
3	20.03	Phthalic acid, but-3-yn-2-yl cyclohexylmethyl ester	C <sub>19</sub> H <sub>22</sub> O <sub>4</sub>	0.42
4	20.89	Ethanone, 1-(7-hydroxy-5-methoxy-2,2-dimethyl-2H-1-benzopyran-8-yl)	C <sub>14</sub> H <sub>16</sub> O <sub>4</sub>	1.79
5	22.69	O-Vanadic acid, tris(4-pentenyl) ester	C <sub>15</sub> H <sub>27</sub> O <sub>4</sub> V	0.19
6	22.78	Androstan-17-ol	C <sub>19</sub> H <sub>32</sub> O	0.13
7	23.12	Pregnane-3,11,20-trione, (5.beta.)	C <sub>21</sub> H <sub>30</sub> O <sub>3</sub>	0.34
8	23.22	2,6,9-Undecatriene	C <sub>11</sub> H <sub>18</sub> O <sub>3</sub>	0.39
9	23.31	2-Pyrimidinamine, 5-bromo-4-methoxy-6-methyl	C <sub>6</sub> H <sub>8</sub> BrN <sub>3</sub> O	0.54
10	23.43	Epiandrosterone	C <sub>19</sub> H <sub>30</sub> O <sub>2</sub>	0.27
11	23.54	Benzene, hexa(1-propenyl)	C <sub>24</sub> H <sub>30</sub>	1.42
12	23.68	Acridin-9-yl-(5-methyl-[1,3,4]thiadiazol-2-yl)-amine	C <sub>16</sub> H <sub>12</sub> N <sub>4</sub> S	0.32
13	23.74	Propanenitrile, 3-[1-[3-(1-pyrrolidinyl)propynyl]-1-cyclohexyloxy]	C <sub>16</sub> H <sub>24</sub> N <sub>2</sub> O	0.3
14	23.86	Lupulon	C <sub>26</sub> H <sub>38</sub> O <sub>4</sub>	47.98
15	24.03	(7-tert-Butyl-2-cyclopropyl-5,6,7,8-tetrahydrobenzo[4,5]thieno[2,3-d]pyrimidin-4-ylsulfanyl)acetic acid	C <sub>19</sub> H <sub>24</sub> N <sub>2</sub> O <sub>2</sub> S <sub>2</sub>	1.2
16	48.55	Phenol, 2,4-bis(1-phenylethyl)	C <sub>22</sub> H <sub>22</sub> O	9.5
17	24.41	Imidazo[1,5-a]pyrrolo[2,1-c][1,4]benzodiazepine-1-carboxylic acid, 7-chloro-9-oxo-11,12,13,13a-tetrahydro, ethyl ester	C <sub>17</sub> H <sub>16</sub> ClN <sub>3</sub> O <sub>3</sub>	1.37
18	24.46	Benz[a]anthracene-7,12-diol, 7,12-dihydro-7,12-dimethyl-, cis	C <sub>20</sub> H <sub>18</sub> O <sub>2</sub>	0.63
19	24.49	Dronabinol	C <sub>21</sub> H <sub>30</sub> O <sub>2</sub>	0.66
20	24.87	26,27-Dinoreergosta-5,22-dien-3-ol, (3.beta.,22E)	C <sub>26</sub> H <sub>42</sub> O	1.95
21	25.00	Phenol, 2,4-bis(1-methyl-1-phenylethyl)	C <sub>24</sub> H <sub>26</sub> O	2.93
22	25.05	3-(1-Acetoxyethyl)-3-phenylpiperidin-2,6-dione	C <sub>15</sub> H <sub>17</sub> NO <sub>4</sub>	1.92
23	25.16	Cholest-20(22)-ene-3,6-dione, (5.alpha.)	C <sub>27</sub> H <sub>42</sub> O <sub>2</sub>	1.81
24	25.20	Tetralin-1-methylamine, N-cyclohexyl-N-oxide	C <sub>17</sub> H <sub>23</sub> NO	0.7
25	25.39	1H-Pyrano[3',4':6,7]indolizino[1,2-b]quinoline-3,14(4H,12H)-dione, 4-ethyl-4-hydroxy-, (S)	C <sub>20</sub> H <sub>16</sub> N <sub>2</sub> O <sub>4</sub>	1.82
26	25.45	Ergosta-7,22-dien-3-ol, (3.beta.,5.alpha.,22Z)	C <sub>28</sub> H <sub>46</sub> O	1.33
27	25.50	Isoindole-1,3(2H)-dione, 5,6-dichloro-2-phenyl	C <sub>14</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>2</sub>	1.11
28	25.60	Androstane-17-carboxylic acid, 3-(acetyloxy)-14-hydroxy-, methyl ester, (3.beta.,5.beta.,14.beta.,17.beta.)	C <sub>23</sub> H <sub>36</sub> O <sub>5</sub>	1.11
29	25.70	D:A-Friedooleanane	C <sub>30</sub> H <sub>52</sub>	1.91
30	25.86	Retinoic acid, 5,6-epoxy-5,6-dihydro-, methyl ester	C <sub>21</sub> H <sub>30</sub> O <sub>3</sub>	0.55
31	25.92	5.alpha.-Androstane, 3.alpha., 11.alpha., 17.beta.-trihydroxy	C <sub>19</sub> H <sub>32</sub> O <sub>3</sub>	0.89
32	26.01	Cholesta-8,14-dien-3-ol, 4,4-dimethyl-, (3.beta.,5.alpha.)	C <sub>29</sub> H <sub>48</sub> O	2.17
33	26.10	Oxalic acid, momoamide, N-(2-chlorophenyl)-, butyl ester	C <sub>12</sub> H <sub>14</sub> ClNO <sub>3</sub>	0.49
34	26.14	N-(2-Amino-4-chlorophenyl)anthranilic acid	C <sub>13</sub> H <sub>11</sub> ClN <sub>2</sub> O <sub>2</sub>	0.37
35	26.21	4H-Pyrido[4,3-b]indole-4,4-dicarboxylic acid, 1,2,3,5-tetrahydro-2-methyl-, diethyl ester	C <sub>18</sub> H <sub>22</sub> N <sub>2</sub> O <sub>4</sub>	1.38
36	26.39	Disilane, 1,1,1,2,2-pentamethyl-2-[(cyclopropyl)(phenylthio)methyl]	C <sub>15</sub> H <sub>26</sub> SSi <sub>2</sub>	0.58

37	26.53	4-(4-Methoxy-6-methyl-5,6,7,8-tetrahydro-[1,3]dioxolo[4,5-g]isoquinolin-5-yl)-5-propyl-2,4-dihydro-pyrazol-3-one	C <sub>18</sub> H <sub>23</sub> N <sub>3</sub> O <sub>4</sub>	0.46
38	26.68	13-Docosenamide, (Z)	C <sub>22</sub> H <sub>43</sub> NO	2.35
39	26.86	1-(1,1,2,3,3,3-Hexafluoropropyl)-4,5-bis(methoxycarbonyl)-1,2,3-triazole	C <sub>9</sub> H <sub>7</sub> F <sub>6</sub> N <sub>3</sub> O <sub>4</sub>	0.7
40	26.91	3-Morpholino-thioacrylomorpholide	C <sub>11</sub> H <sub>18</sub> N <sub>2</sub> O <sub>2</sub> S	0.17
41	27.02	2-(4,7-Dimethyl-quinazolin-2-ylamino)-6-methyl-pyrimidin-4-ol	C <sub>15</sub> H <sub>15</sub> N <sub>5</sub> O	0.21
42	27.23	Estra-1,3,5(10)-trien-17-one, 3,7-dihydroxy-, O-methyloxime, (7.alpha.)	C <sub>19</sub> H <sub>25</sub> NO <sub>3</sub>	0.15
43	27.35	Androstane, (5.beta.,14.beta.)	C <sub>19</sub> H <sub>32</sub>	0.27
44	28.19	Benzeneacetic acid, 3-methoxy-4-[(trimethylsilyl)oxy]-, ethyl ester	C <sub>14</sub> H <sub>22</sub> O <sub>4</sub> Si	0.25
45	28.85	Cyclodeca[b]furan-2(3H)-one, 9-(acetyloxy)-3a,4,5,8,9,11a-hexahydro-4-hydroxy-6,10-dimethyl-3-methylene	C <sub>17</sub> H <sub>22</sub> O <sub>5</sub>	0.17
46	29.37	5,5',8,8'-Tetrahydroxy-3,3'-dimethyl-2,2'-binaphthalene-1,1',4,4'-tetrone	C <sub>22</sub> H <sub>14</sub> O <sub>8</sub>	0.29
47	29.60	Benzenamine, 4,4',4''-phosphinyldynetris[N,N-dimethyl	C <sub>24</sub> H <sub>30</sub> N <sub>3</sub> OP	0.19
48	29.69	Quinoline, 2-chloro-6-methoxy-4-methyl	C <sub>11</sub> H <sub>10</sub> ClNO	0.15

Table S2. Analytical results of serum samples by the proposed method.

Sample	Without spiking (mM)	Glucose spiked (mM)	Glucose measured (mM)	Recovery (%)
Serum -1	3.28 ± 0.27	10	12.82 ± 0.79	95.40
Serum -2	3.06 ± 0.55	20	23.91 ± 1.12	104.25
Serum -3	3.58 ± 0.81	50	54.89 ± 1.63	102.62
Serum -4	3.28 ± 0.44	100	107.33 ± 1.54	104.13