Supplementary Table. 1 The Alog *P*, hydrogen bond number, and total polarity surface of benzenoids from *Antrodia cinnamomea* determined by molecular modeling

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
| Compound | Molecular weight (Da) | Alog *P* | Hydrogen-bond acceptor number | Hydrogen-bond donor number | Total polarity surface |
| Compound 1 | 184.19  | 1.799  | 4 | 2 | 58.92  |
| Compound 2 | 198.22  | 2.025  | 4 | 1 | 47.92  |
| Compound 3 (CoQ0) | 182.17  | 0.510  | 4 | 0 | 52.60  |
| Compound 4 | 154.16  | 1.815  | 3 | 2 | 49.69  |
| Compound 5 | 168.15  | 1.600  | 4 | 2 | 58.92  |
| Compound 6 | 196.20  | 2.051  | 4 | 0 | 36.92  |