Table S1 Pivotal structural parameters of the compounds M1-M10.

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| Parameters Optimized Active site Parameters Optimized Active site  |
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| --- | --- | --- | --- | --- | --- | --- |
| M1 |  |  | M2 |  |  |  |
| C6-O7 | 1.2221 | 1.2220 | C3-C17 | 1.5103 | 1.5105 |
| C12-O17 | 1.3413 | 1.3414 | C17-O18 | 0.4195 | 0.4196 |
| O17-H25 | 0.9871 | 0.9873 | O18-H28 | 0.9626 | 0.9631 |
| O19-H27 | 0.9875 | 0.9875 | C1-O20-H30 | 107.05 | 107.05 |
| O18-H26 | 0.966 | 0.9660 | C12-O19-H29 | 107.1 | 107.08 |
| C11-C12-O17-H25 | 180.0 | 0.4009 | C3-C17-O18-H28 | 180.0 | 43.19 |
| C2-C1-O19-H27 | 180.0 | 128.31 |  |  |  |  |
| C2-C3-O18-H26 | 0.0 | -157.34 |  |  |  |  |
| M3 |  |  | M4 |  |  |  |
| C17-O19 | 1.2099 | 1.2103 | C14-O1 | 1.365 | 1.3652 |  |
| C17-O18 | 1.3437 | 1.3441 | O1-C21 | 1.4393 | 1.4399 |  |
| O18-H28 | 0.9706 | 0.9699 | O5-H30 | 0.9662 | 0.966 |  |
| C7-O15-H27 | 107.19 | 107.19 | O3-H29 | 0.9919 | 0.9923 |  |
| C14-O21-H29 | 107.38 | 107.36 | C19-O5-H30 | 110.20 | 110.26 |  |
| C9-C17-O18-H28 | 180.0 | -109.84 | C13-O3-H29 | 105.91 | 105.88 |  |
|  |  |  | O8-C14-O1-C21 | -84.43 | 177.53 |  |
| M5 |  |  | M6 |  |  |  |
| C6-O13 | 1.3408 | 1.3411 | C17-O2 | 1.3702 | 1.3701 |  |
| O13-H22 | 0.9852 | 0.9863 | O2-C15 | 1.4249 | 1.425 |  |
| C16-O19 | 1.3409 | 1.3412 | O8-H50 | 0.9662 | 0.9654 |  |
| O19-H29 | 0.9855 | 0.985 | O10-H51 | 0.9921 | 0.9925 |  |
| C2-C6-O13-H22 | 0 | -52.46 | C18-C17-O2-C15 | 63 | 56.17 |  |
| C10-C16-O19-H29 | 0 | -48.52 | C25-C26-O8-H50 | -179.98 | -168.85 |  |
|  |  |  | C22-C27-O10-H51 | 0.5401 | 23.83 |  |
| M7 |  |  | M8 |  |  |  |
| C12-O18 | 1.364 | 1.3639 | O2-C17 | 1.3772 | 1.3779 |  |
| O18-C26 | 1.4129 | 1.4126 | C20-O7 | 1.2128 | 1.2118 |  |
| O17-H37 | 0.9857 | 0.9858 | C27-O9 | 1.3529 | 1.3521 |  |
| O16-H36 | 0.9857 | 0.9859 | O9-H50 | 0.9639 | 0.9635 |  |
| C11-C12-O18-C26 | 16.23 | -116.95 | C31-O10 | 1.4208 | 1.4213 |  |
| C9-C14-O17-H37 | -0.084 | -36.02 | O10-H51 | 0.9632 | 0.9626 |  |
| C4-C5-O16-H36 | -0.139 | 0.2294 | C18-C17-O2-C15 | 109.41 | 137.82 |  |
|  |  |  | C21-C27-O9-H50 | -178.4 | -10.29 |  |
|  |  |  | C29-C31-O10-H51 | -66.81 | 112.95 |  |
| M9 |  |  | M10 |  |  |  |
| C18-O2 | 1.3567 | 1.3584 | C16-O2 | 1.3619 | 1.3625 |  |
| C16-O2 | 1.391 | 1.3905 | C14-O2 | 1.4053 | 1.4057 |  |
| C21-O7 | 1.1862 | 1.186 | O5-H40 | 0.9719 | 0.9729 |  |
| C28-O9 | 1.3379 | 1.3381 | O3-H38 | 0.9641 | 0.9636 |  |
| O9-H49 | 0.9408 | 0.9407 | O9-H50 | 0.9637 | 0.9641 |  |
| C32-O11 | 1.1806 | 1.1805 | C26-O9 | 1.3518 | 1.3518 |  |
| C32-O10 | 1.3266 | 1.3266 | C17-C16-O2-C14 | -141.44 | 111.44 |  |
| O10-H50 | 0.9459 | 0.9465 | C20-C26-O9-H50 | -176.91 | 35.54 |  |
| C19-C18-O2-C16 | 101.97 | 87.28 |  |  |  |  |
| C23-C28-09-H49 | -177.57 | -177.32 |  |  |  |  |
| C30-C32-O10-H50 | 179.99 | -75.41 |  |  |  |  |

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Table S2 Electron densities and Laplacian values of electron densities for M1 at both status

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| 化学键 M1 ρbcp e/A-3 DockM1ρbcp e/A-3 M1 ∇2ρbcpe/A-5 DockM1∇2ρbcpe/A-5 |
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| --- | --- | --- | --- | --- |
| C1-C2 | 2.123 | 1.997 | -21.194 | -20.926 |
| C1-O19 | 2.085 | 1.895 | -9.610 | -8.473 |
| O19-H27 | 2.273 | 2.084 | -47.384 | -31.848 |
| C2-H21 | 1.905 | 1.719 | -24.158 | -17.534 |
| C2-C3 | 2.146 | 2.019 | -21.371 | -20.898 |
| C3-O18 | 1.994 | 1.819 | -7.283 | -7.208 |
| O18-H26 | 2.462 | 2.234 | -50.385 | -37.323 |
| C3-C4 | 2.097 | 1.976 | -20.959 | -20.479 |
| C4-H22 | 1.947 | 1.717 | -25.849 | -17.443 |
| C4-C5 | 2.159 | 2.048 | -21.251 | -21.709 |
| C5-C16 | 2.022 | 1.914 | -18.819 | -19.430 |
| C16-C14 | 1.941 | 1.814 | -18.068 | -18.145 |
| C14-O15 | 2.471 | 2.315 | -2.364 | -10.725 |
| C13-C14 | 1.919 | 1.791 | -17.751 | -17.832 |
| C8-C13 | 2.024 | 1.913 | -18.862 | -19.457 |
| C5-C6 | 1.813 | 1.679 | -16.174 | -15.721 |
| C6-O7 | 2.703 | 2.513 | 4.952 | -6.037 |
| C6-C8 | 1.823 | 1.690 | -16.301 | -15.936 |
| C12-C13 | 2.061 | 1.929 | -19.746 | -19.383 |
| C11-C12 | 2.117 | 1.985 | -21.285 | -20.522 |
| C12-O17 | 2.072 | 1.869 | -9.445 | -7.617 |
| O17-H25 | 2.278 | 2.118 | -47.451 | -29.614 |
| C8-C9 | 2.149 | 2.039 | -21.165 | -21.636 |
| C9-H23 | 1.948 | 1.736 | -25.637 | -17.882 |
| C9-C10 | 2.074 | 1.956 | -19.981 | -20.348 |
| C10-C11 | 2.139 | 2.020 | -20.933 | -21.312 |
| C11-H24 | 1.925 | 1.703 | -24.824 | -17.182 |
| C10-C20 | 1.728 | 1.607 | -14.798 | -14.858 |
| C20-H28 | 1.891 | 1.667 | -23.516 | -16.626 |
| C20-H29 | 1.867 | 1.669 | -22.913 | -16.653 |
| C20-H30 | 1.867 | 1.670 | -22.911 | -16.711 |

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Table S3 Electron densities and Laplacian values of electron densities for M6 at both status

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| 化学键 M6 ρbcp e/A-3 DockM6ρbcp e/A-3 M6 ∇2ρbcpe/A-5 DockM6∇2ρbcpe/A-5 |
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| --- | --- | --- | --- | --- |
| C17-C21 | 1.984 | 2.012 | -20.620 | -21.120 |
| C17-O2 | 1.691 | 1.740 | -6.437 | -6.727 |
| O2-C15 | 1.399 | 1.519 | -3.771 | -4.775 |
| C21-H42 | 1.785 | 1.709 | -18.562 | -17.276 |
| C21-C26 | 2.009 | 2.015 | -20.888 | -20.984 |
| C26-O8 | 1.739 | 1.823 | -6.834 | -7.188 |
| O8-H50 | 2.065 | 2.240 | -28.475 | -37.430 |
| C25-C26 | 2.004 | 2.006 | -20.906 | -20.929 |
| C25-H44 | 1.823 | 1.735 | -19.453 | -17.869 |
| C25-C19 | 2.039 | 2.035 | -21.581 | -21.483 |
| C18-C19 | 1.900 | 1.899 | -19.246 | -19.188 |
| C18-C20 | 1.756 | 1.731 | -17.130 | -16.589 |
| C20-O7 | 2.224 | 2.392 | -9.986 | -8.380 |
| C20-C22 | 1.796 | 1.760 | -17.561 | -16.989 |
| C22-C23 | 1.928 | 1.927 | -19.783 | -19.751 |
| C19-C24 | 1.672 | 1.666 | -15.635 | -15.494 |
| C24-O9 | 2.401 | 2.509 | -8.998 | -6.274 |
| C23-C24 | 1.737 | 1.714 | -16.790 | -16.349 |
| C22-C27 | 1.922 | 1.928 | -19.513 | -19.534 |
| C27-C30 | 1.967 | 1.987 | -20.442 | -20.791 |
| C27-O10 | 1.857 | 1.895 | -8.096 | -8.297 |
| O10-H51 | 1.794 | 2.045 | -22.393 | -33.317 |
| C23-C28 | 2.037 | 2.031 | -21.500 | -21.388 |
| C28-H45 | 1.810 | 1.734 | -19.187 | -17.857 |
| C28-C29 | 1.945 | 1.956 | -20.307 | -20.419 |
| C29-C30 | 2.030 | 2.030 | -21.514 | -21.576 |
| C30-H46 | 1.804 | 1.725 | -19.006 | -17.662 |
| C29-C31 | 1.583 | 1.609 | -14.326 | -14.866 |
| C31-H49 | 1.756 | 1.670 | -18.139 | -16.713 |
| C31-H47 | 1.734 | 1.670 | -17.757 | -16.669 |
| C31-H48 | 1.733 | 1.668 | -17.747 | -16.649 |

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