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|-------------------|-------------------|---------------------------|------------------------|--------------------------|---------------------|-------------------------|
| 1. MW | 5. max Hbonds | 9. atomic density | 13. potential Hbonds | 17. DBD correlation | 21. ligand distance | 25. C3 distance |
| 2. charge | 6. polarity | 10. bb entropy loss | 14. frustration index | 18. LBD correlation | 22. centrality | 26. C4 distance |
| 3. hydrophobicity | 7. polarizability | 11. sc entropy loss | 15. bfactor | 19. residue correlations | 23. C1 distance | 27. sequence separation |
| 4. aromaticity | 8. flexibility | 12. solvent accessibility | 16. structural entropy | 20. DNA distance | 24. C2 distance | |

