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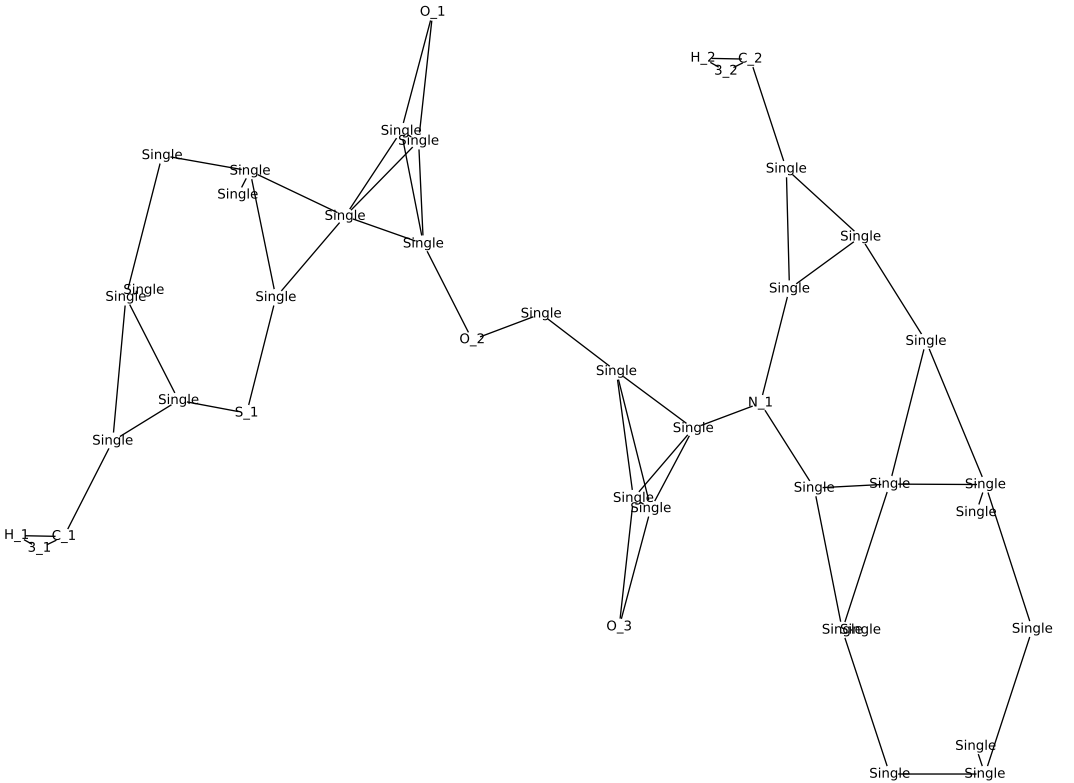
O_1
H_2
C_2
3_2
Single
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