

Fig. S1 PADs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3)\rightarrow CO_2+H$ at E_c of 8.6 kcal mol⁻¹ from QCT calculations on the YMS PES.

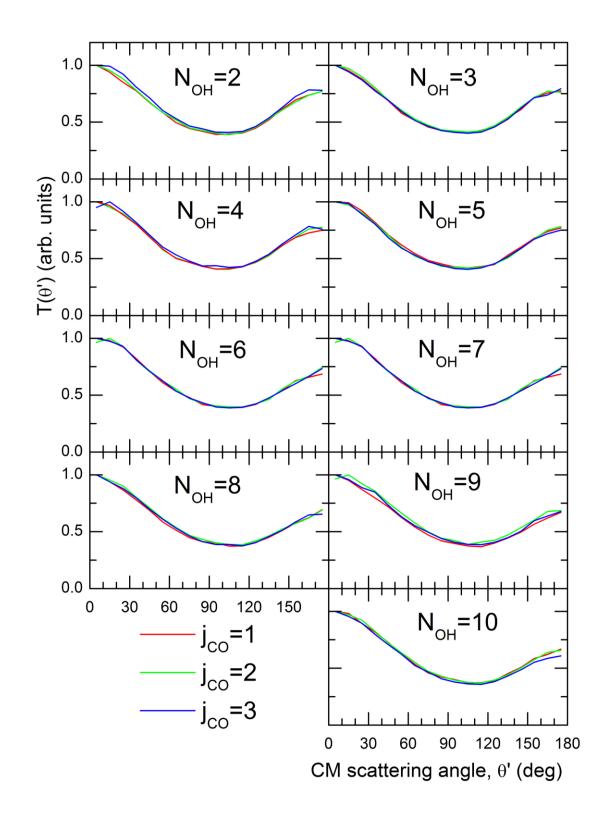


Fig. S2 PADs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3)\rightarrow CO_2+H$ at E_c of 14.1 kcal mol⁻¹ from QCT calculations on the YMS PES.

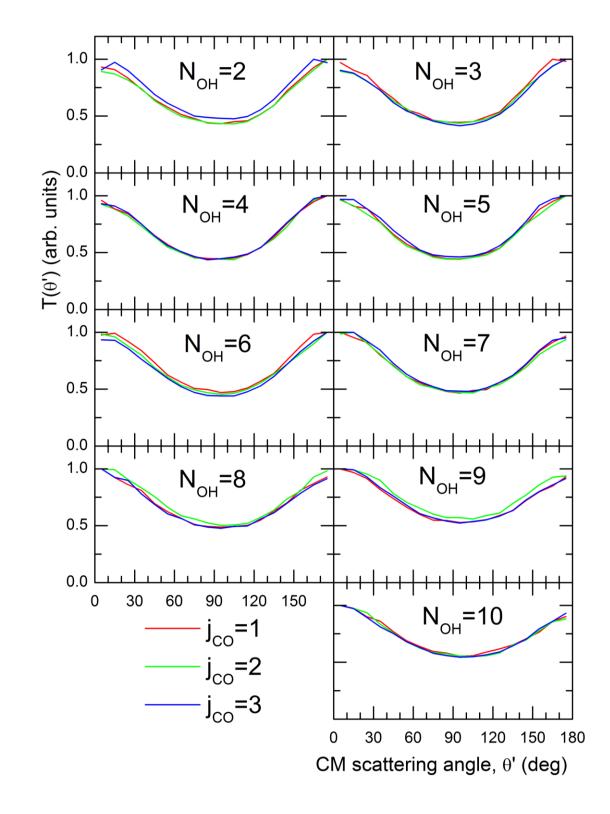


Fig. S3 PADs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3) \rightarrow CO_2+H$ at E_c of 8.6 kcal mol⁻¹ from QCT calculations on the LTSH PES.

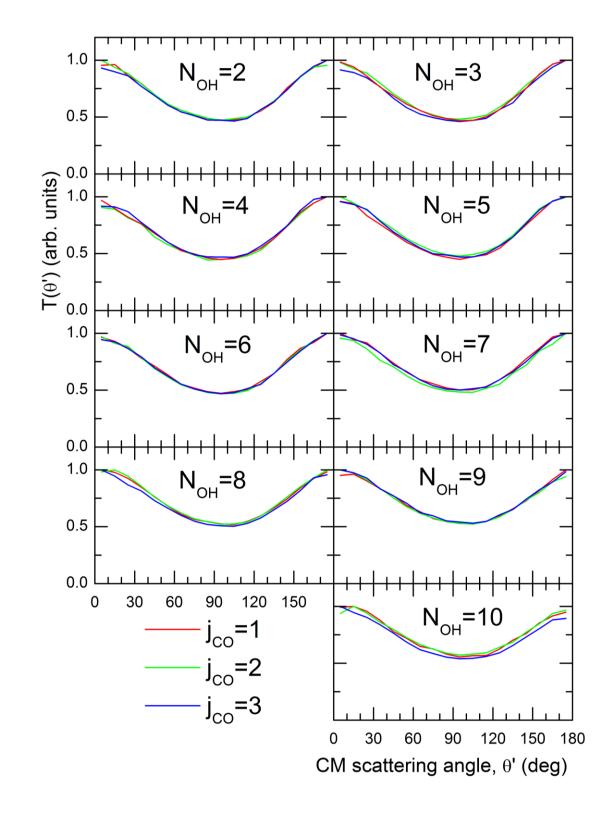


Fig. S4PADs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3)\rightarrow CO_2+H$ at E_c of 14.1 kcal mol⁻¹ from QCT calculations on the LTSH PES.

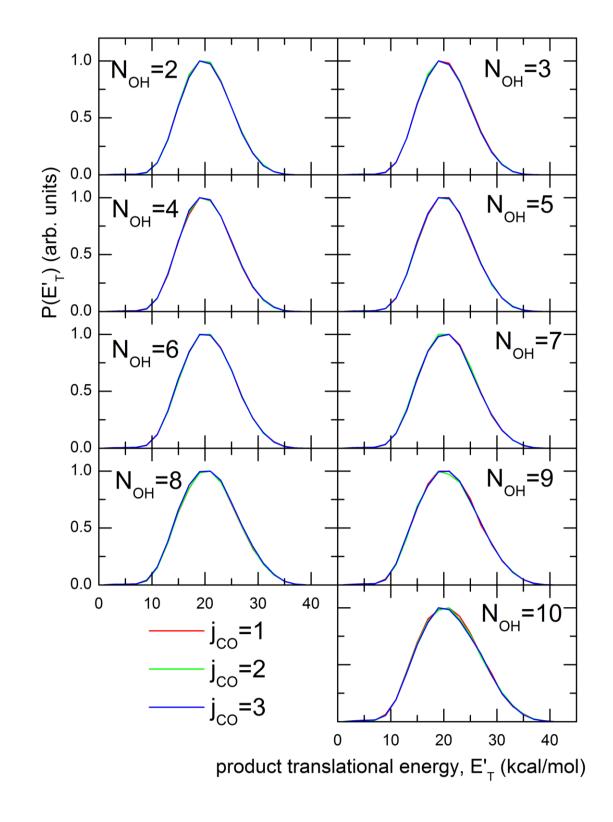


Fig. S5 PTDs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3) \rightarrow CO_2+H$ at E_c of 8.6 kcal mol⁻¹ from QCT calculations on the YMS PES.

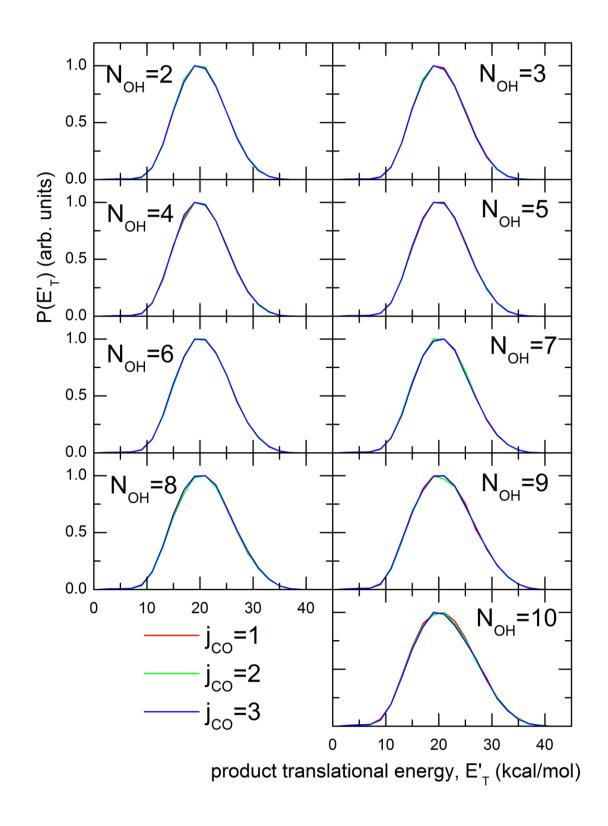


Fig. S6 PTDs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3)\rightarrow CO_2+H$ at E_c of 14.1 kcal mol⁻¹ from QCT calculations on the YMS PES.

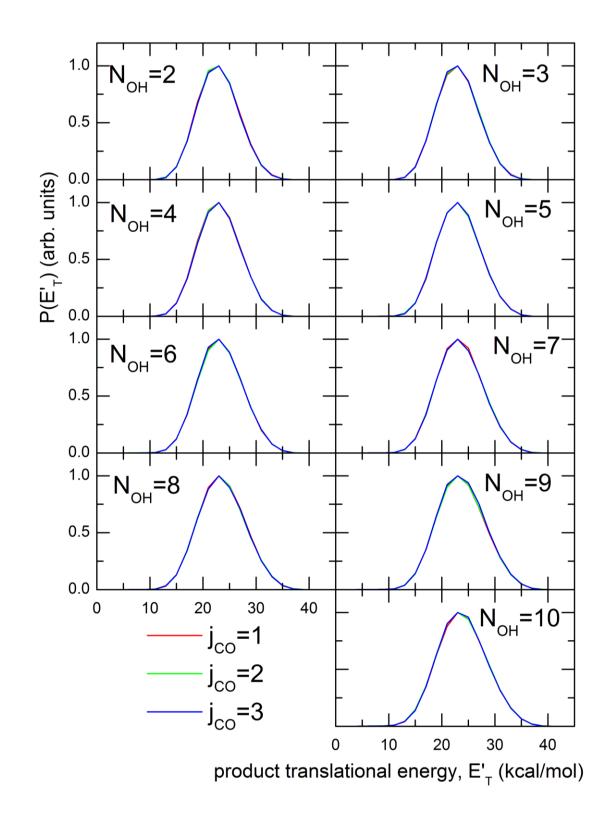


Fig. S7 PTDs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3) \rightarrow CO_2+H$ at E_c of 8.6 kcal mol⁻¹ from QCT calculations on the LTSH PES.

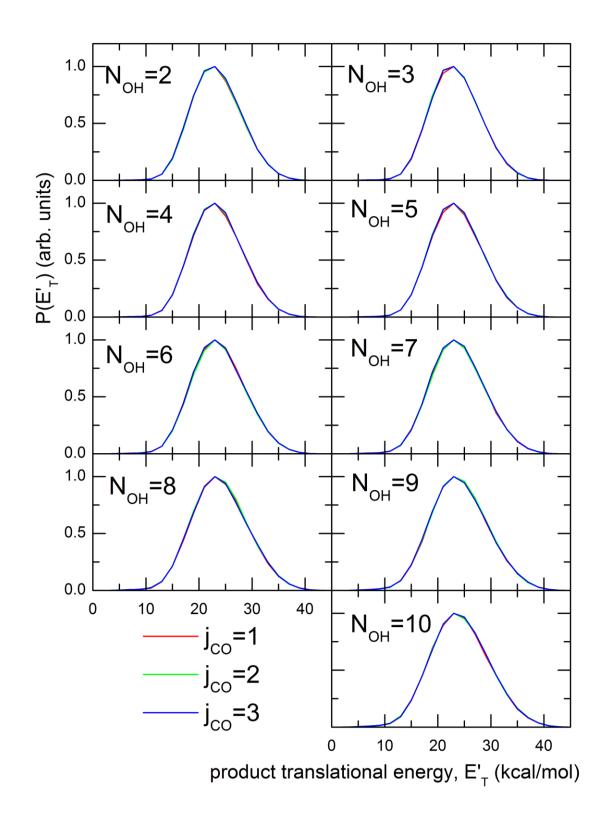


Fig. S8 PTDs for different OH rotational levels for the reactions $OH(N)+CO(j=1,2,3)\rightarrow CO_2+H$ at E_c of 14.1 kcal mol⁻¹ from QCT calculations on the LTSH PES.