

Abstract Submitted  
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**Energetics of reactions in a dielectric barrier discharge with argon carrier gas: Halocarbons** SEAN WATSON, Polytechnique Montreal, BERNARD NISOL, Ghent University, STEPHAN REUTER, MICHAEL R. WERTHEIMER, Polytechnique Montreal — The novel method we developed for understanding energy exchanges between argon (Ar) carrier gas and precursor molecules in a large-area  $216\text{ cm}^2$  dielectric barrier discharge (DBD) reactor has resulted in a series of articles, each relating to a different family of organic compounds. This communication focuses on two new groups, perfluorocarbons,  $C_xF_y$  and perchlorocarbons,  $C_xCl_y$ , and compares results with earlier ones for hydrocarbons,  $C_xH_y$ [1] and hydrofluoromethanes,  $CH_xF_y$ [2]. The precursors (in  $\text{o}/\text{o}$  concentrations) were mixed with Ar in a 20 kHz, 8 kV (peak to peak) DBD. For each separate compound  $E_m$ , the energy (in eV) absorbed per molecule, was determined from measurements of the time resolved discharge current,  $I_d$ , and the gap voltage,  $V_{gap}$ . Optical emission spectra were recorded and  $E_m$  was plotted as a function of precursor flow rate,  $F_d$ . The process generally led to thin plasma polymer (PP) deposits (e.g. on Si wafer substrates). Their characteristics, like their C/F or C/Cl composition ratios from XPS measurements, strongly correlated with  $E_m$  and  $F_d$ , as did PP deposition rates and water contact angles. [1] B. Nisol et al., Plasma Process Polym, 2016;14:e201600191 [2] S. Watson et al., Plasma Process Polym, 2020;17:e201900125.

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