## Shape optimization for LDOS inside a cativity

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## Adjoint method for shape optimization

The local density of states (LDOS) is proportional to the imaginary party of the dyadic green's function [2]:

$$\rho(\mathbf{x}_0, \boldsymbol{\omega}) = \frac{\boldsymbol{\omega}}{\pi c^2} Im Tr[\overline{\overline{G^{EP}}}(\mathbf{x}_0, \mathbf{x}_0, \boldsymbol{\omega}) + \overline{\overline{G^{HM}}}(\mathbf{x}_0, \mathbf{x}_0, \boldsymbol{\omega})]$$
(1)

We only consider the electric LDOS  $\rho_e$  which comes from the electric part of the green's function  $\overline{\overline{G^{EP}}}(\mathbf{x}_0,\mathbf{x}_0,\boldsymbol{\omega})$ . The electric field  $\mathbf{E}(\mathbf{x}_0,\boldsymbol{\omega})$  is determined by the green's function by:

$$\mathbf{E}(\mathbf{x}_0, \boldsymbol{\omega}) = i\mu_0 \boldsymbol{\omega} \int \overline{\overline{G^{EP}}}(\mathbf{x}_0, \mathbf{x}, \boldsymbol{\omega}) \cdot \mathbf{j}(\mathbf{x}) d^3 \mathbf{x}$$
 (2)

In the time harmonic case with only electric dipole at  $\mathbf{x_0}$  as the source, we have  $\mathbf{j}(\mathbf{x_0}) = \frac{\partial \mathbf{p}(\mathbf{x_0})}{\partial t} = -i\omega\mathbf{P}\delta^3(\mathbf{x} - \mathbf{x_0})$ , inserting this into Eq. (2) we have

$$\mathbf{E}(\mathbf{x}_0, \boldsymbol{\omega}) = \frac{\boldsymbol{\omega}^2}{\varepsilon_0 c^2} \overline{\overline{G^{EP}}}(\mathbf{x}_0, \mathbf{x}_0, \boldsymbol{\omega}) \cdot \mathbf{P}$$
(3)

Compare to the LDOS expression Eq. (1), we have

$$\rho_e(\mathbf{x_0}) = \frac{\varepsilon_0}{\pi \omega} \frac{1}{|\mathbf{P}_0|} Im \sum_j \hat{\mathbf{s}}_j \cdot \mathbf{E}_{s_j}(\mathbf{x}_0)$$
(4)

where  $\mathbf{E}_{s_j}$  denotes the field from a dipole source at  $\mathbf{x}_0$  polarized in the  $\mathbf{s}_j$  direction, with an unit dipole moment  $\mathbf{P}_0 = \mathbf{s}_j$  (thus  $|\mathbf{P}_0| = 1$ ), and the sum over j accounts for all possible orientations. This is a small modification to Eq. (8) in Owen's paper [1], where we make the electric dipole to be unit dipole and the coefficient  $\varepsilon_0$  to the front.

Therefore, we can get the LDOS of a specified by three scattering simulation.

In Owen's thesis, we have that for a shape deformation  $\delta x_n(\mathbf{x}')$ , the variation for the object function F is (Eq. (5.28) in Ref. [3]).

$$\delta F = 2Re \int \delta x_n(\mathbf{x}') [(\varepsilon_2 - \varepsilon_1) \mathbf{E}_{\parallel}(\mathbf{x}') \cdot \mathbf{E}_{\parallel}^{\mathbf{A}}(\mathbf{x}') + (\frac{1}{\varepsilon_1} - \frac{1}{\varepsilon_2}) \mathbf{D}_{\perp}(\mathbf{x}') \cdot \mathbf{D}_{\perp}^{\mathbf{A}}(\mathbf{x}')] dA$$
 (5)

However, this equation is wrong because it assumed that  $\partial F/\partial \mathbf{E}$  is the complex conjugate of  $\partial F/\partial \bar{\mathbf{E}}$ . In fact,

$$\overline{\left(\frac{F}{\partial \mathbf{E}}\right)} = \frac{\partial \bar{F}}{\partial \bar{\mathbf{E}}} \neq \frac{\partial F}{\partial \bar{\mathbf{E}}} \tag{6}$$

Therefore, the right variation for the object function should be

$$\delta F = \int \delta x_n(\mathbf{x}') [(\varepsilon_2 - \varepsilon_1) \mathbf{E}_{\parallel}(\mathbf{x}') \cdot \mathbf{E}_{\parallel}^{\mathbf{A}}(\mathbf{x}') + (\frac{1}{\varepsilon_1} - \frac{1}{\varepsilon_2}) \mathbf{D}_{\perp}(\mathbf{x}') \cdot \mathbf{D}_{\perp}^{\mathbf{A}}(\mathbf{x}')] dA + \text{Conjugate Adjoint.}$$
(7)

Since the electric LDOS  $\rho_e$  is the imaginary part of some function, say  $\tilde{\rho}_e$ . The optimum of  $\tilde{\rho}_e$  must also be the optimum of  $\rho_e$ . So we can first take the object function to be  $\tilde{\rho}_e$ , for each direction j in Eq. (4), the source of the adjoint field is a electric dipole at  $\mathbf{x}_0$  with the amplitude

$$\frac{\partial \tilde{\rho_e}}{\partial \mathbf{E}_{s_i}} = \frac{\varepsilon_0}{\pi \omega} \int dx^3 \delta(\mathbf{x} - \mathbf{x}_0) \hat{\mathbf{s}_j}$$
 (8)

while the conjugate adjoint field part  $\frac{\partial \tilde{p}_e}{\partial \tilde{\mathbf{E}}_{s_i}} = 0$ .

This is a also a dipole at  $\mathbf{x} = \mathbf{x}_0$  in the  $\hat{\mathbf{s}}_j$  direction, so the adjoint field is the same as the original field up to a constant scalar:

$$\mathbf{E}_{s_j}^{\mathbf{A}}(\mathbf{x}) = \frac{\varepsilon_0}{\pi \omega} \mathbf{E}_{sj} \tag{9}$$

Therefore, we have the variation for the electric LDOS as

$$\delta \rho_e = \frac{\varepsilon_0}{\pi \omega} Im \sum_{j} \int \delta x_n(\mathbf{x}') \{ (\varepsilon_2 - \varepsilon_1) [\mathbf{E}_{s_j \parallel}(\mathbf{x}')]^2 + (\frac{1}{\varepsilon_1} - \frac{1}{\varepsilon_2}) [\mathbf{D}_{s_j \perp}(\mathbf{x}')]^2 \} dA$$
 (10)

In scuff-scatter, we can use --EPFile to get the electric field of specified points, however, it only applies to points away from the scattering surface. For points on the surface, the obtained data are just garbage. Instead, we can implement the option --PSDFile to obtain the effective electric and magnetic surface charge and current. In the scuff calculation, the effective electric and magnetic surface charge and current are defined as

$$\sigma = \mathbf{n} \cdot \mathbf{D} \tag{11a}$$

$$\mathbf{K} = \mathbf{n} \times \mathbf{H} \tag{11b}$$

$$\eta = \mathbf{n} \cdot \mathbf{B} \tag{11c}$$

$$\mathbf{N} = \mathbf{n} \times \mathbf{E} \tag{11d}$$

Therefore, the parallel electrical field and perpendicular displacement can be computed as

$$\mathbf{E}_{\parallel} = \mathbf{N} \tag{12a}$$

$$\mathbf{D}_{\perp} = \sigma \mathbf{n} \tag{12b}$$

After the surface fields are obtained, we can first verify this adjoint method on a void sphere case. For a void sphere, the electromagnetic surface modes satisfy:

$$\varepsilon_m(E)H_l(k_m a)[k_d a J_l(k_d a)]' = \varepsilon_d J_l(k_d a)[k_m a H_l(k_m a)]'$$
(13)

where a corresponds to the void radius, l is the (integer) index denoting the angular momentum,  $k_m = \sqrt{\varepsilon_m} k_0$  and  $k_d = \sqrt{\varepsilon_d} k_0$  are wave vectors in metal and void.  $J_l$  and  $H_l$  are spherical Bessel and Hankel functions of the first kind[4, 5].

By solving Eq. (13), we can quickly get some resonant parameters and use it as the initial guess for the general shape optimization purpose.

Next Step:

• Modify the runMesher code and make spherical harmonic basis

## References

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