# Efficient 3-D Extraction of Interconnect Capacitance Considering Floating Metal Fills With Boundary Element Method

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Abstract-Inserting dummy (area fill) metals is necessary to reduce the pattern-dependent variation of dielectric thickness in the chemical-mechanical polishing (CMP) process. Such floating dummy metals affect interconnect capacitance and, therefore, signal delay and crosstalk significantly. To take the floating dummies into account, an efficient method for three-dimensional (3-D) capacitance extraction based on boundary element method is proposed. By introducing a floating condition into the direct boundary integral equation (BIE) and adopting an efficient preconditioning technique, and the quasi-multiple medium (QMM) acceleration, the method achieves very high computational speed. For some typical structures of area fill, the presented algorithm has shown over  $1000 \times$  speedup over the industry-standard Raphael while preserving high accuracy. Compared with the recently proposed PASCAL in the work of Park et al. (2000), the proposed method also has about ten times speedup. Since the dummies are not regarded as normal electrodes in capacitance extraction, the proposed method is much more efficient than the conventional method, especially in cases with a large number of floating dummies.

*Index Terms*—Boundary-element method (BEM), design for manufacturability, dummy fill, interconnect capacitance extraction.

### I. INTRODUCTION

**F** OR current very-large-scale integration (VLSI) circuit, chemical-mechanical polishing (CMP) is a necessary manufacturing step by which the wafer is polished with a rotating pad and slurry to achieve the planarized surfaces [1], [2]. Because the dielectric thickness strongly depends on the pattern density of underlying metal layer, a widely used method for reducing the variation of dielectric thickness in the CMP process is to insert dummy metals (this procedure is also called "area fill") [1]–[4]. These dummies are situated between signal lines to increase the pattern density and, at the same time, influence the electrical characteristics of interconnects in different ways depending on whether they are grounded or on the floating state. In the application-specific integrated circuit (ASIC) design, the floating dummy fills are preferred due to the short design period and considerable area to be filled.

The authors are with the Department of Computer Science and Technology, Tsinghua University, Beijing 100084, China (e-mail: yu-wj@tsinghua.edu.cn). Digital Object Identifier 10.1109/TCAD.2005.853690 capacitance and, therefore, signal delay and crosstalk [1]–[3]. Nowadays, the area fill synthesis considering the impact on

Such floating dummy fills have a strong impact on interconnect

circuit performance has become an important problem of the design for manufacturability (DFM) [4]. For the electrical characterization or optimization of design rules for the dummy fills, a huge number of simulations are required for the structures involving floating dummy fills. However, most of the available parasitic extraction tools do not have specific treatment for floating dummies. For the case with floating conductors, a conventional field solver (such as SYNOPSYS Raphael) treats them as normal electrodes and extracts the full capacitance matrix. Then, the capacitance matrix is reduced by considering that the total charge on each floating conductor is equal to zero [5]. Obviously, the cost of central processing unit (CPU) time for extracting the full capacitance matrix is prohibitive while including a lot of floating dummies.

On the other hand, three-dimensional (3-D) interconnect capacitance extraction with the boundary element method (BEM) has drawn much attention recently, because of its advantage of dimensionality reduction over the finite difference method and finite element method. Many fast algorithms based on BEM have been proposed, including the fast multipole approach [6], hierarchical approach [7], the precorrected fast Fourier transform (FFT) algorithm [8], and the quasi-multiple medium (QMM) method [9], [10]. Different from other kinds of BEM, the QMM-accelerated BEM employs the direct boundary integral equation (BIE) [13], which has the property of resulting in a sparse coefficient matrix for a multiregion problem. By exploiting this character to enlarge the matrix sparsity and efficient techniques of equation organization and solution, the QMM-BEM solver has shown ten times speedup and memory saving over the multipole-accelerated method (FastCap 2.0) [6], for 3-D capacitance extraction [10].

In this paper, the QMM-accelerated BEM is extended to efficiently handle the structure with floating dummy fills. An equation corresponding to the floating condition is added into the discretized BEM equations, for each floating conductor. Then, with a new preconditioning technique for iterative equation solution, the interconnect capacitances are directly obtained. Since the floating conductor is not considered as normal electrode and no redundant capacitance entries need to be solved and reduced, the proposed method is much more efficient than the conventional method, particularly in the case of a great number of dummy fills in modern ASIC design.

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Some typical structures of area fill are calculated to demonstrate the efficiency of the proposed method. Compared with Raphael (RC3), which is the finite difference solver with advanced nonuniform meshing, the proposed method has a speedup ratio above several hundreds.

In [11] and [15], two algorithms of capacitance extraction were proposed with specific consideration of the floating dummy fills. The algorithm in [15] is based on a so-called fictitious domain method, which uses the Lagrange multiplier  $\lambda$  to consider the conductor boundary and replaces the complex shape domain of potential computation with a simple one. However, no other algorithm was compared with the algorithm based on fictitious domain method, and the computational results in [15] did not show high efficiency (it costs 1 h for a structure including 130 floatings). The basic idea of [11] is similar to the authors, but the algorithm (called PASCAL) is implemented based on the finite difference method. Further work on electrical characterization of metal fills using the PASCAL can be found in [12]. Since the authors are not able to obtain the PASCAL for direct comparison in the same computational environment, an indirect comparison is carried out based on the speedup ratios to Raphael. For similar structures, the proposed method finally shows about ten times speedup over PASCAL.

The remainder of this paper is organized as follows. In Section II, the QMM-based capacitance extraction for structure without floating conductors is briefly reviewed. Efficient techniques handling the structure with floating dummies are presented in Section III. Experimental results are reported in Section IV, and conclusions are drawn in Section V.

## II. THREE-DIMENSIONAL CAPACITANCE EXTRACTION WITH THE QMM-ACCELERATED BEM

In direct BEM, the Laplace equation fulfilled by the electrical potential u in each homogenous dielectric region can be transformed into the following direct BIEs [9], [10], [13]

$$cu + \int_{\Gamma_i} q^* u d\Gamma = \int_{\Gamma_i} u^* q d\Gamma \qquad i = 1, \dots, M \qquad (1)$$

where c is a constant depending on the boundary geometry and  $\Gamma_i$  is the boundary of dielectric region *i* (assuming there are totally *M* regions). *q* is the normal electrical field intensity and  $q = \partial u / \partial n$  where *n* means the vector outward normal to boundary. For 3-D space, the fundamental solution  $u^*$  is  $1/4\pi r$ and  $q^* = \partial u^* / \partial n$ . Employing constant quadrilateral elements to make boundary discretization and evaluating the direct BIE at collocation points (one for each element), the discretized BIEs are obtained from (1) for each dielectric region. After evaluating two types of boundary integral, a set of linear equations with unknowns of *u* and *q* is obtained [9].

Moreover, u and q fulfill the compatibility equations along the interface of two adjacent dielectrics a and b

$$\begin{cases} \varepsilon_{a} \frac{\partial u_{a}}{\partial n_{a}} = -\varepsilon_{b} \frac{\partial u_{b}}{\partial n_{b}} \\ u_{a} = u_{b} \end{cases}$$
(2)



Fig. 1. Typical 3-D interconnect capacitor with five dielectric layers is cut into a  $3 \times 2$  structure.

where  $\varepsilon_a$  and  $\varepsilon_b$  stand for the permittivities of dielectrics a and b, respectively. With (2) the *u* and *q* unknowns on the same interfacial element for both adjacent dielectric regions can be united, respectively. Therefore, the discretized BIEs for all dielectric regions can be put together. By substituting the boundary conditions (*u* is known on conductor surfaces as bias voltage, and *q* is supposed to be zero on the Neumann boundary as shown in Fig. 1), and preserving only the items of unknown variables to the left side of the equal sign, an overall linear system is obtained

$$\mathbf{A}\boldsymbol{x} = \boldsymbol{f} \tag{3}$$

where x is a vector comprising all discretized unknowns of u and q. Theoretically, any arrangement of the discretized BIEs in (3) is correct. However, without careful consideration, the population of the coefficient matrix A would be too chaotic to solve the equation efficiently. In [9], Yu *et al.* present an effective arrangement of the unknowns and collocation points, as well as the storage scheme for the resulting sparse matrix A, which facilitates the matrix–vector multiplication in GMRES [14] equation solution for problem involving a large number of regions.

The above process can be regarded as the conventional direct BEM for a multiregion problem. A localization character is revealed by formula (1), where the variables in each BIE are within the same dielectric region. This character of direct BEM results in a blocked sparse coefficient matrix A for a multidielectric problem. For example, a typical capacitor with three dielectrics is shown in Fig. 2, and the corresponding sparse matrix A is shown in Fig. 3 (the effective equation arrangement is used). In the QMM method, every actual dielectric is decomposed into some fictitious medium blocks, whose permittivities are all the same as the original dielectric, to increase the sparsity of matrix A. With the storage technique of sparse blocked matrix and efficient iterative equation solver such as the preconditioned GMRES algorithm, the computing time and memory usage for the original problem will be greatly reduced. In practice, each layer of an actual multilayered interconnect structure is decomposed into  $m \times n$  fictitious medium blocks, perpendicular to the bottom substrate plane (see Fig. 1). In order to decrease the additional efforts brought by the QMM decomposition, the cutting planes are also dispersed along the x-axis or y-axis uniformly.



Fig. 2. Structure with three dielectrics (cross-sectional view).



Fig. 3. Matrix population corresponding to the structure in Fig. 2 and the regions with dashed-line contour represent the added entries when considering the floating conductors.

In [10], an efficient approach to automatically determine the QMM cutting was proposed to maximize the sparsity of the overall coefficient matrix. Moreover, an efficient preconditioning technique was proposed for the GMRES solution, which resulted in about 30% less solution time than that using the Jacobi preconditioner. Finally, the enhanced QMM–BEM solver showed over ten times speedup and memory saving over the multipole approach with comparable accuracy [10].

## III. EFFICIENT TECHNIQUES CONSIDERING THE FLOATING DUMMIES

In this section, first, the basic idea of the proposed method is given, which introduces the floating condition. The details of equation organization and preconditioning are then discussed. Finally, the proposed method is compared with the straightforward method for handling floating dummies.

## A. Basic Idea

For a structure including  $N_c$  interconnect conductors, an approach of setting the *j*th conductor to 1 V and the rest to 0 V is usually used to determine the self and coupling capacitances of the *j*th conductor (it is also called master conductor). This procedure can be repeated  $N_c$  times to get the capacitance matrix, which includes all self and coupling capacitances among the interconnects. If there is no floating dummy in the structure, the potential u is known on all conductor surfaces. This boundary condition is utilized to form (3) in Section II. After solving (3), the self-capacitances and coupling capacitances can be evaluated by the integral of the normal electrical field intensity q on the conductor surfaces [6]–[10].

If some conductors in the simulated structure are changed to be floating (without setting known voltage), (3) cannot be solved because there are more unknowns (potential u on these floating conductors) than equations. So, the task at hand here is to supply additional equations about the floating dummies and make the enlarged system of linear equations solvable. The proposed method of capacitance extraction for the structure involving floating dummies includes six steps.

- 1) Set the *j*th interconnect conductor to 1 V and the rest interconnects to 0.
- For each dielectric region, discretize its boundary (including dummy surface, since it is also a part of boundary) and formulate the discretized BIEs as in Section II.
- Put the discretized BIEs for all regions together with (2), and substitute the known conditions on interconnect conductor surface and the Neumann boundary, as in Section II.
- Supply some equations about the floating dummies so that the total number of equations is equal to the number of unknowns (including the additional *u* unknowns of dummies).
- 5) Solve the generated system of linear equations and get the charges on interconnect conductors that equal the capacitances.
- 6) Repeat steps 1)–5) with different voltage settings to get all capacitances among the interconnects.

Now, methods on how to supply the equations about the floating dummies are discussed. The following equation is used to calculate the total charge of a floating dummy

$$\int_{\Gamma_f} \sigma \mathrm{d}\Gamma = \int_{\Gamma_f} \varepsilon q \mathrm{d}\Gamma = \overline{Q}_f \tag{4}$$

where  $\Gamma_f$  is the surface of the dummy conductor, and  $\sigma$  is the surface charge density.  $\varepsilon$  is the permittivity of the dielectric surrounding surface  $\Gamma_f$ , and q is the normal electrical field intensity. Here, the charge  $\overline{Q}_f$  is usually zero for each dummy as its initial electrical state. Since the dummy surface is discretized into elements, (4) is actually used as its discretized form, which involves discretized q unknowns on a dummy surface.

If there are  $N_f$  floating dummies,  $N_f$  new equations of (4) are derived, one for each dummy. On the other hand, as has been discussed, there are also  $N_f$  additional u unknowns of dummies, one for each dummy (because of the equipotential property). Hence, the extended equation system from (3) can be solved and the interconnect capacitances are then obtained.

#### B. Equation Organization and Preconditioning

As discussed above, more unknowns and equations are introduced for the case involving floating dummy fills. They are organized together with the initial discretized BEM equations in (3) to form a new overall linear system. Putting (4) and the unknowns of potential on floating dummies at the end, the new linear system is derived with the form

$$\begin{bmatrix} \mathbf{A} & \mathbf{C1} \\ \mathbf{C2} & \mathbf{0} \end{bmatrix} \cdot \begin{bmatrix} x \\ u_f \end{bmatrix} = \begin{bmatrix} f \\ \overline{q} \end{bmatrix}, \quad \text{i.e., } \mathbf{A}' x' = f' \quad (5)$$

where **A**, x, and f are obtained from the linear system (3) where the floating character of dummies is not considered.  $u_f$  is the vector consisting of the unknown potential of dummies, and  $\bar{q}$  is made up of the right-hand side of (4) (actually a zero vector in the context of this paper). **C1** and **C2** are corresponding submatrices of coefficients. For the problem with three dielectrics in Fig. 2, if some conductors are changed to be floating, the modification of the overall coefficient matrix can be illustrated as the regions with dashed-line contour in Fig. 3.

Below, details of calculating the submatrices C1 and C2 with discretizing the direct boundary integrals (1) and (4) are given for a problem involving multiple dielectric regions and multiple floating dummies. C1 corresponds to the coefficients of the new unknowns  $u_f$  and can be expressed as

$$(\mathbf{C1})_{ij} = \int_{\partial\Omega_{f,j}^{(i)}} q_i^* \mathrm{d}\Gamma$$
 (6)

where  $q_{(i)}^*$  means the normal derivative of  $u^*$  in (1), which is related with the *i*th collocation point.  $\partial \Omega_{f,j}^{(i)}$  represents the part of surface of the *j*th floating dummy, which is within the same dielectric region with the collocation point *i*. **C2** corresponds to the coefficients in (4), and can be expressed as

$$(\mathbf{C2})_{ij} = \begin{cases} \int_{\Gamma_j} \varepsilon d\Gamma, & \text{when column } j \text{ corresponds} \\ & \text{to a } q \text{ unknown on dummy } i \\ 0, & \text{otherwise} \end{cases}$$
(7)

where  $\Gamma_j$  stands for the *j*th boundary element, and  $\varepsilon$  is the permittivity of the surrounding dielectric.

For capacitance extraction without floating dummies, a socalled extended Jacobi (EJ) preconditioner for GMRES algorithm was proposed in [10]. It is an approximation to  $A^{-1}$ , and makes the convergence rate at least 30% faster than using the Jacobi preconditioner, while introducing little overhead [10]. However, since there is a zero diagonal block in A', the EJ preconditioner cannot be used here directly. Assume an EJ preconditioning matrix **P** is constructed for the original coefficient matrix **A**, then the new preconditioner is constructed as follows:

$$\mathbf{P}' = \begin{bmatrix} \mathbf{P} & \mathbf{0} \\ \mathbf{0} & \mathbf{I} \end{bmatrix} \tag{8}$$

where I stands for the identity matrix. Since the dimension of I (equal to the number of dummies) is much smaller than the total number of discretized unknowns,  $\mathbf{P}'$  still approximates  $(\mathbf{A}')^{-1}$  to some extent. Therefore, the new preconditioner



Fig. 4. Straightforward method of capacitance extraction for structure involving floating dummies. (a) Network of capacitances to be extracted. (b) Desired capacitance network after eliminating the floating nodes.

should also improve the rate of convergence for the GMRE equation solver. Numerical experiments verified this analysis.

#### C. Comparison With the Straightforward Method

The straightforward method in conventional field solvers treats the floating dummies as normal electrodes, extracts the capacitances related with these dummies, and reduces them to get capacitances of interconnect. To illustrate this, a problem involving three interconnect conductors and two floating dummies is taken for example (see Fig. 4). If the coupling capacitances between conductor 1 and other interconnect conductors (2 and 3) ought to be calculated, i.e., conductor 1 is master conductor, all capacitances related with floating dummies [f1 and f2 in Fig. 4(a)] are first calculated. This invokes computing twice, with a floating dummy set to 1 V and the rest to 0 V. Then, conductor 1 is set to 1 V to calculate its capacitances. After that the capacitance network shown in Fig. 4(a) is obtained. Since f1 and f2 are floating, the network can be reduced to the one shown in Fig. 4(b) by eliminating the nodes of f1 and f2. Finally, the equivalent capacitance between conductor 1 and conductor 2 (or 3) is obtained, as if dummies do not exist. In fact, these equivalent capacitances are what the authors are concerned about.

For capacitance extraction of a structure involving  $N_c$  interconnect (signal) conductors and  $N_f$  floating dummies, the straightforward method is compared with the proposed method as follows.

1) To calculate the capacitances related with one master conductor (usually for the analysis of critical path), a 3-D field solver must be invoked for  $(N_f + 1)$  times in the straightforward method, and then network reduction is performed to eliminate the floating nodes. In contrast, the 3-D field solver is invoked only once in the proposed method, although the degree of the linear system to be solved increases for  $N_f$  due to the additional unknowns of dummy potential. Usually in the field solution, the number of discretized unknowns is much larger than  $N_f$ . So the increased computational expense of the proposed method in the once field solution is very little. Ignoring both the increased expense of the proposed method and the time of network reduction in straightforward method, the speedup of the proposed method can be expected to be  $(N_f + 1)$  for capacitance extraction with one master conductor.



Fig. 5. Two conductor lines surrounded with 24 square floating dummies located at two layers.

2) To calculate the whole capacitance matrix of the interconnects, the 3-D field solver needs to be invoked for  $(N_f + N_c)$  times in the straightforward method, while in the proposed method only  $N_c$  times are needed. Similarly, the proposed method has about  $(N_c + N_f)/N_c$  times speedup over the straightforward one when extracting the whole capacitance matrix. This is also remarkable since  $N_f$  is usually much larger than  $N_c$ .

Therefore, for both the one-master extraction and the wholematrix extraction, the proposed method has large speedup over the straightforward one. When there are a lot of dummies ( $N_f$ or  $N_f/N_c$  is large), the proposed method is particularly efficient compared with the straightforward method.

#### **IV. NUMERICAL RESULTS**

The proposed algorithm has been implemented in QBEM [10], a software prototype for 3-D capacitance extraction using the QMM-accelerated BEM. Some typical structures of floating dummy fills have been calculated to verify the efficiency of the proposed algorithm. All experiments are run on a Sun Ultra V880 with frequency of 750 MHz.

Three examples are shown in Figs. 5-7, respectively. The first one involves two parallel conductor lines surrounded with 24 square floating dummies located at two layers. The size of each conductor lines is  $1 \times 23 \times 0.6$ ; the size of each dummy metal is  $1 \times 1 \times 0.6$ , and they all are within a window of size  $9 \times 23 \times 4$ , above the grounded substrate. The second one is a typical structure of floating dummy fills, which are arranged as a dot array in the oblique alignment and is often adopted for minimizing changes in the coupling capacitances and maximizing the uniformity of the pattern density [2], [11]. The size of each parallel conductor line is  $6 \times 0.5 \times 0.5$ ; the size of each dummy metal is  $0.5 \times 0.5 \times 0.5$ , and they are all within a window of size  $6 \times 9 \times 1.5$ . In example 3 (Fig. 7), the dummy insertion rule is similar to that in the second example (Fig. 6), as dot array in the oblique alignment. Different from example 2, the conductor lines and dummies are located at two layers and the former constructs a  $1 \times 2$  crossover. In this example, two conductor lines in the same layer are of size  $6.5 \times 0.5 \times 0.5$ , while the third one is of  $0.5 \times 6.25 \times 0.5$ . The size of each dummy is  $0.5 \times 0.5 \times 0.5$ , and all are within



Fig. 6. Three conductor lines and 34 floating dummies of the dot-array type.



Fig. 7. Three conductor lines and 53 floating dummies of the dot-array type located at two layers.

a window of size  $6.5 \times 6.25 \times 2.5$ . All length parameters above are in micrometer. The relative permittivities of layers in the same case are different, with values 1.9, 2.9, and 3.9.

In each of the three structures, conductor 1 is set to be a master (see Figs. 5–7), and its self-capacitance and coupling capacitances with other conductors are computed using Raphael and the proposed method. In the computation with Raphael, the dummy-related capacitances are calculated by setting each dummy as master conductor one after another, and the wanted capacitances are outputted after performing circuit reduction [5]. In the computations with the proposed method,  $2 \times 1$ ,  $1 \times 3$ , and  $4 \times 2$  QMM cutting are applied, respectively, for the three structures to achieve high speed. The capacitance results and other detailed data for these three examples are listed in Table I.

In Table I, it can be seen that the errors of capacitances computed with the proposed method are all within 3% (using Raphael's result as criterion), while the speedup ratio to Raphael ranges from 300 to several thousands. The proposed method uses less than one-tenth of the memory used by Raphael. It should also be pointed out that the electrical potential of dummy metal is also calculated, and it is very close to Raphael's result as well (with discrepancy within 5%). The number of dummies and number of unknowns in the proposed method are also listed in Table I, noting that the latter is larger than the element number because each interfacial element has two unknowns. From these data, it can be seen that the dummy number is much larger than the number of signal conductors, while the number of discretized unknowns is much larger than the number of dummies. They support some assumption for analysis in Section III.

TABLE I COMPARISON OF COMPUTATIONAL RESULTS WITH OUR METHOD AND RAPHAEL (CAPACITANCE IN  $10^{-18}$  F)

		Dummy N	C <sub>11</sub>	C <sub>12</sub>	C <sub>13</sub>	Ele N	Var N	Iter	Time (second)	Memory (megabyte)	Speed-Up
First	Proposed Method	24	2446	-1091		654	792	37	1.1	1.7	299
Example	Raphael		2470	-1088		$85 \times 10^{3*}$			328.4	18	
Second	Proposed Method	24	766.1	-11.4	-10.4	1006	1172	55	1.9	2.7	1330
Example	Raphael	34	753.2	-11.5	-11.5	$281 \times 10^{3*}$			2527	54	
Third	Proposed Method	50	732.1	-161.1	-141.5	1789	2501	52	3.0	5.9	2311
Example	Raphael	33	741.6	-163.0	-161.4	$385 \times 10^{3*}$			6932	72	

Dummy\_N means the number of dummies.

Ele\_N means the number of elements in our method.

Var\_N means the number of unknowns.

Iter means the number of iterative steps.

\* Number of default grids generated by Raphael automatically.

 TABLE II

 Comparison of Computational Time of the Proposed Method and the Conventional Method for Example 1

	Time for Extractin	g Capacitances H	Related With (second)	Total Time	Speed-Up Ratio for	Speed-Up Ratio for	
	Conductor 1	Conductor 2	24 Dummios	(second)	One-Master	Whole-Matrix Extraction	
	Conductor	Conductor 2	24 Dummes		Extraction		
Conventional	0.84	0.84	25.07	26.75	1	1	
Method	0.01	0.01	20.07	20.75	1	1	
Proposed Method	1 1.13	1.13		2.26	22.9	11.8	

To demonstrate the efficiency of the new preconditioner proposed in Section III-B, the problem is also solved without a preconditioner and with a modified Jacobi preconditioner. Experiment results show that without preconditioner the GMRES algorithm cannot converge within 300 steps for the last two problems and converges with 230 steps for the first problem. While using the modified Jacobi preconditioner, the iterative numbers for the three examples are 39, 62, and 62, respectively. Here, the modified Jacobi preconditioner is constructed also as (8). So, it is clear that the modified EJ preconditioner performs very well for the capacitance extraction involving floating dummies.

To demonstrate the advantage of our method over the conventional one, the whole capacitance matrix of example 1 is calculated using both methods. With the proposed method, QBEM ran twice, setting the two signal lines as master, respectively. With the conventional method, QBEM without the presented modification ran totally 26 times setting all conductors including dummies as master: one for each time. The details of computational time are listed in Table II. Without considering the time to eliminate the floating nodes, the speedup ratio of the proposed method is found to be 22.9 and 11.8, for single-master extraction and whole-matrix extraction, respectively. This is consistent with the theoretic analysis in Section III-C, which expects the speedup ratio to be 25 and 13. Because no more source of error is induced by the proposed method, it has the same accuracy with the conventional one.

In [11], a fast method to extract interconnect capacitance considering the floating dummy fills was implemented as software PASCAL, whose performance was also compared with Raphael. They used several cases of dummies located as dot arrays in the oblique alignment (similar to those in Fig. 6 and Fig. 7) to make a comparison, which showed that PASCAL has a speedup ratio varying from 11 to 290 (see [11, Table 2]) as the number of dummy metals increases. Since the PASCAL for direct comparison in the same computational environment



Fig. 8. Speedup ratio of PASCAL and the proposed method to Raphael for structures involving different numbers of dummies.

cannot be obtained, only a relative comparison is carried out based on the information provided by [11]. Fig. 8 shows the speedup ratios of two methods to Raphael for similar structures involving oblique alignment of dot dummies. From this, it can be expected that the method in this paper is about ten times as fast as the PASCAL.

#### V. CONCLUSION

Based on the QMM-accelerated BEM [9], [10], an efficient method is proposed for the interconnect capacitance extraction involving dummy fills. By introducing the floating condition directly, only unknowns with quantity of dummy metals are added to the linear system corresponding to the same extraction problem except that the dummies are not floating. Since no redundant capacitances related with the dummies need to be solved and reduced, the proposed method is particularly efficient in the case of a great number of floating dummies. With the quasi-multiple medium technology and efficient preconditioned GMRES solver, the proposed method finally shows several thousand times speedup to Raphael while preserving high accuracy, for typical filling structures. Besides, the proposed method also shows about ten times speedup to the fast PASCAL in [11] for similar structures.

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