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USE OF FIELD INFORMATION TO MATCH THE RECORDS IN TWO FILES

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Consider two files of records. Within each file, each record corresponds to a different population unit; but the two files correspond to the same general population. We want to identify "matches," i.e., pairs of records (from the two files) that each correspond to the same population unit.

Each record contains data in K fields which correspond to characteristics such as age, race, etc. We may observe patterns of agreement/disagreement among the fields, for each pair of records. Using this information, we want as best as possible to identify matches. The problem of how best to use the field information has been addressed for K=3, under assumption that the events "agreement in field i," i=1, ..., K are stochastically mutually independent -- for true matches and likewise for true nonmatches. We address the problem for K>3, and avoid reliance on the assumption of independence by fitting interaction terms which reflect stochastic positive dependences.

Key words: record linkage, field dependence, convex programming, log-linear model.

1. Problem Definition. Suppose we have two files "A" and "B" containing N $_{\mbox{\scriptsize A}}$ records and N $_{\mbox{\scriptsize B}}$ records respectively. Each A-record corresponds to a different population unit, likewise each B-record. The two files correspond to the same general population; and our goal is to identify "linkages": a linkage is a coupling of an A-record and a B-record that correspond to the same population unit. In other words, we want to decide which pairings correspond to the same population unit, and which do not. We refer to Fellegi and Sunter (1969) for a lengthy discussion of many aspects of this general, frequently arising situation, including their "method of moments," which provides a starting point for the method of this paper. (Because of the directions that this paper takes, and the resulting complications, we depart from the notation of that paper. disrespect is intended.) As that paper indicates, these files may represent samples from larger populations. We presume that, by whatever means available, we have "unduplicated" each file so that, in essence, no two records within a file correspond to the same population unit. Coulter (1985) suggests one scheme for this "unduplication."

As in our reference list, the Coulter paper appeared in the Proceedings of a May 1985 workshop on exact matching methodologies, co-sponsored by the Internal Revenue Service, the Washington Statistical Society, and the Federal Committee on Statistical Methodology. All our references on record linkage appeared in these lengthy proceedings and, in some cases, appeared earlier elsewhere.

There are $N_A N_B$ possible pairings of A-records and B-records. Out of these, the maximum possible number of "true" matches (same population unit) is $\min(N_A, N_B)$. Hence we introduce $F_O = \min(N_A, N_B)/N_A N_B$ (1)

as the maximum possible fraction of pairings that are true matches; we do not know how many pairings are true matches, but

 F_0 gives an upper bound. Typically, however, our files, and the population that underlies them, will be "blocked" into disjoint subsets in a parallel fashion so as to guarantee that, if a population unit appears in both files A and B, it will with virtual certainty appear in the same block for both. (If not, we will not be able to identify the true match for this population unit.) Kelley (1985) and Winkler (1985) discuss strategies for blocking; later in this paper we discuss the subject. For now, suppose that the files contain N_{hA} and N_{hB} records in block h. We consider the possible pairings of A and B within each block. Summing over all blocks, we get $\sum N_{hA}N_{hB}$ and $\sum \min(N_{hA},N_{hB})$ as number of possible pairs and maximum possible number of true matches. As maximum possible fraction we now have

$$F_{O} = \left[\sum_{h} \min(N_{hA}, N_{hB})\right] / \left(\sum_{h} N_{hA} N_{hB}\right) . \tag{2}$$

Typically, F_0 will be close to 0. (If external information permits us to determine an upper bound which is even smaller, we could use that in place of (2). The actual fraction may in practice be considerably less than F_0 . For this reason one ought to carry out analyses of this paper using a smaller F_0 , as a sensitivity check.)

For each pair we compare K data fields that appear in both the A- and B-records; these fields typically correspond to such characteristics as age, race, or street name. There are 2^K cells corresponding to patterns of field agreement/disagreement; each pair falls into one and only one of the cells. We will number the fields from 1 to K. Let S denote a subset of these fields. We designate "cell S" as that cell for which there is agreement for the fields in S, and disagreement for the other fields. Note the terminology: agree/disagree pertains to fields, and match/nonmatch to population units. Thus cell S corresponds to the vector outcome γ , associated with a pair, that appears in many papers on record linkage. As an example let K=5 and $S = \{2,3,5\}$. The cell S consists of those pairs for which there

is agreement in fields 2, 3 and 5, and disagreement in fields 1 and 4.

Let P_S denote the observed proportion of all pairs that land in cell S. Let F denote the fraction of all pairs that are true matches. Let Q_{1S} denote the proportion of all true match pairs that fall into cell S, and Q_{2S} denote the same for true nonmatches. We have

$$P_S = FQ_{1S} + (1-F)Q_{2S}$$
 (3)

Although we observe P_S , we cannot observe its components F, Q_{1S} and Q_{2S} ; we only know $F \le F_0$. Using only the known P_S and F_0 , we will fit values for F, Q_{1S} and Q_{2S} .

Once these values are fitted, Fellegi and Sunter (in their Section 2) use Q_{1S} and Q_{2S} to form a rule for deciding whether a pair represents a true match. The procedure corresponds to a hypothesis test in its formation of critical regions based on largest and smallest values of the likelihood ratio - here, Q_{1S}/Q_{2S} - and in its consideration of Type I and Type II error probabilities. Hence this is one major use of Q_{1S} and Q_{2S} . Henceforth we let

U_S denote Q_{1S}/Q_{2S} .

Instead of the hypothesis test, in this paper we use Q_{1S} and Q_{2S} to develop a posterior probability of true match: that is, the probability of a true match, given that a pair is in cell S. From this point one can make an assignment of file-A records to file-B records by solving a transportation-type linear-programming problem so as to maximize the posterior expected number of true matches, i.e., the sum of posterior probabilities of match, in the assignment. But instead, we might simply specify a cutoff decision rule that a posterior probability exceeding, say, .90, corresponds to true match (and less than .90 does not). Interpretation is simplified by the fact that posterior probabilities largely appear to gravitate toward either

0 or 1. We are inclined to favor investigation of the cutoff decision rule, and thus will not deal here with the intricacies of making an assignment based on the transportation problem.

Hence we have discussed uses of Q_{1S} and Q_{2S} once they are fitted. For the most part this paper considers how best to fit Q_{1S} and Q_{2S} , in advance of using them. We will consider fits which give an end result of form

log
$$U_S = \text{constant plus} \sum_{i \in S} w_i$$
 (4)

with $w_i>0$; this model appears in many of the Proceedings papers on record linkage. A (relative) weight w_i is being assigned to field i. By using (4) we maintain a proper hierarchy among likelihood ratios: if (the set of agreeing fields for) S is strictly contained in (same for) cell T, we want $U_S < U_T$. In such an instance, T represents a clear-cut advance beyond S with respect to field agreements; and we want $U_S < U_T$ so that for T the choice "match" is favored over the choice "non-match" more than it is for S. This hierarchy must be maintained; it is not enough merely to fit Q_{1S} and Q_{2S} which are positive.

The relationship in (4) is maintained under the "independence" model which we present at the start of the next section. Such a model has been used in other papers on record linkage. However, we do not wish to operate under this assumption of independence among field agreements. (This assumption states that, for true-match pairs, the events "agreement in field i" and "agreement in field j" are, for i # j, stochastically independent, and that for true-nonmatch pairs the same thing holds. But typically, for match or nonmatch or both, there will be positive dependence between these events for at least some (i,j). Rather than consider the possible robustness of results based on the independence assumption against positive dependence, we want to obtain results which take the positive dependence into account.) Thus we will develop a log-linear model which, as far as possible, accounts for positive dependences among field agreements - while preserving the

relationship in (4), and the accompanying hierarchy for \mathbf{U}_{S} . The development and use of this latter model is the primary idea of the paper.

In regard to our model and method, we stress that the only needed inputs are the observable cell proportions P_S . We do not rely on a sample of pairs which are known to be true matches (or true nonmatches) (e.g., Tepping (1968)) for information. (With F_O small, true matches are very scanty.) Nor do we make use of the transmission rates of Howe and Lindsay (1981): probabilities (known) that a recorded field entry is the true entry.

We will use the following example as illustration. We considered files for (A) the 1980 Decennial Census, and (B) the Post-Enumeration Survey for that census. Each record corresponds to a person. We use K=5 fields:

- (1) A combination of race and ethnicity
- (2) A combination of gender and marital status
- (3) Date of birth (month and year)
- (4) Surname
- (5) A combination of street name and location within community.

We considered only those records for which all fields were filled. With "blocking" based on "Census Bureau numbering area," there were 49 blocks. The number of total pairs was 1321100, the number of possible true matches 4465, and the value of F_0 .003380. The 32 cells, with accompanying P_S , are in Table 1; the 5 integers preceding P_S indicate field agreement (1) or field disagreement (0) for the 5 fields. That is, 11000

	Table 1.	Values of PS	
00000 .334568 10000 .353175 01000 .063819 11000 .080692 00100 .000318 10100 .000372 01100 .000129	00010 .000724 10010 .001824 01010 .000121 11010 .000522 00110 .000000 10110 .000017 01110 .00005	00001 .046498 10001 .081199 01001 .010829 11001 .018664 00101 .000098 10101 .000179 01101 .000052	00011 .000751 10011 .002542 01011 .000172 11011 .000709 00111 .00016 10111 .000100 01111 .000202 11111 .001266

indicates agreement on fields 1 and 2, and disagreement for 3-4-5. As we see, slightly more than 2/3 of all pairs disagree on all of fields 2-3-4-5.

2. Outline of Method. This section attempts an overview of the paper. We first address the assumption of independence: consider a model of form

$$Q_{1S} = \begin{bmatrix} \pi_{i} \in S \\ i \in S \end{bmatrix} \begin{bmatrix} \pi_{i} \in S \\ i \in S \end{bmatrix} (1 - y_{i})$$
(5)

We express Q_{2S} likewise with u_i replacing y_i . Let E_i denote the event "agreement in field i." Then y_i corresponds to $P(E_i)$, conditional, given that there is a true match. The model (5) states that the events E_1 , ..., E_K are mutually independent, given that there is a true match. Similar mutual independence holds, given that there is a true non-match, with u_i corresponding to conditional $P(E_i)$. Although we later seek to avoid this assumption of mutual independence, we use the above independence model to develop preliminary results. These initial results, in Sec. 3, will in essence extend the 3-field results of Fellegi and Sunter (their Sec. 3.32 and Appendix 2). From (5), the resulting value of w_i in (4) is seen to be

$$w_i = \log(y_i(1-u_i)/u_i(1-y_i))$$
 (6)

under this independence model. We will fit $u_i < y_i$, so that $w_i > 0$.

Let $\mathbf{p_i}$ denote the fraction of all pairs for which there is agreement in field i, $\mathbf{p_{ij}}$ denote the fraction for which there is agreement in fields i and j, and so forth. As a needed step in developing our method, we must consider

$$R_{ij} = p_{ij} - p_i p_j, \text{ for } i \neq j.$$
 (7)

Incidentally, in our example the values of \boldsymbol{p}_{i} are

For the smallest values of p_i , agreement (for field i) is the rarest - and, plausibly, the corresponding fields i discriminate the most usefully between matches and non-matches.

In practice we will typically have all ${\rm R}_{\mbox{ij}} > 0\,;$ if not, we cannot proceed further. With

$$p_{i} = Fy_{i} + (1-F)u_{i}$$
 (9)

$$p_{ij} = Fy_iy_j + (1-F)u_iu_j$$
 (10)

under independence as in (5) and with $y_i > u_i$ presumed, the measured positivity of R_{ij} is due to a combination of three possible sources:

- (a) Departure of F from O.
- (b) Excess of y_i over u_i and of y_j over u_j .
- (c) Positive dependence between E_i and E_j that the independence model does not reflect (that is, (10) understates the value of p_{ij}).

(To see how (a) and (b) contribute to the positivity of R_{ij} , one substitutes (9) and (10) into (7) and simplifies:

$$R_{ij} = F(1-F)(y_i-u_i)(y_j-u_j).)$$

In Sec. 3 we will fit a value for F, which we denote by F^* , under the assumption of independence. Accompanying this value for F will be a fit of values for y_i and u_i , which we denote by y_i^* and u_i^* , for $i=1,\ldots,K$. As the Sec. 3 formulas (11), (12), and (18) will indicate, these fitted values depend heavily on the values R_{ij} . In effect, we will partition the positivity of the R_{ij} 's into the above three sources (a), (b) and (c), in doing this fit.

The fitted value F^* will typically exceed F_0 , the (logical) maximum possible F. It would be wrong to assign this excess to

effect (a): departure of F from 0. Thus we will presume that any excess of fitted F* over maximal F_0 is, instead, attributable to effect (c): positive dependence. Accordingly, in preparation for Sec. 4 where (corresponding to the value of F_0) we measure positive dependence, at the end of Sec. 3 we replace y_i^* and u_i^* by a pair of values y_i^* and u_i^* which correspond to F_0 . We choose y_i^* and u_i^* so as to hold effect (b) (excess of y over u) constant: we maintain the spacing between y_i^* and u_i^* in the sense that the value of w_i in (6) is preserved.

Then, in Sec. 4 we use F_0 , y_1' and u_1' , and a log-linear model, to fit a set of interaction terms that measure the positive dependence of (c). Having fitted these terms, we adjust the table proportions P_S to compensate for interaction effects. We obtain a new "de-interactionized" table for which, with $F=F_0$, the assumption of independence could more nearly plausibly hold. Accordingly, as in Sec. 5, we fit a new set of values y_1' and u_1' , and fit additional interaction effects. As explained there, we continue to repeat these steps alternately as long as we continue to obtain an improved fit. (We always have $y_1^* > u_1^*$ and $y_1' > u_1'$. But if either $u_1' \le 0$ or $y_1' \ge 1$ is obtained, we cannot proceed further, in these iterations.)

Thus in summary, we fit results under the independence assumption (Sec. 3), fit interaction terms and de-interactionize (Sec. 4), and repeat these two steps (Sec. 5) as long as forward progress is obtained. Then, in Sec. 6, we discuss final results and their interpretation; we also consider different possible values for the true F. Sec. 7 discusses how best to form fields from a given data base, and Sec. 8 discusses computation issues.

3. Results under Independence. This section fits Q_{1S} and Q_{2S} for K fields under the independence assumption — and then makes the adjustment needed for Sec. 4. As a start, consider three fields i, j and k. For 3 fields under the independence assumption Fellegi and Sunter solve 7 equations (corresponding to 2^3-1 cell probabilities P_S) in 7 unknowns (F,y_i,u_i) to obtain fitted F^* given by:

$$C_{ijk} = (R_{ij}R_{ik}R_{jk})^{1/2};$$
 (11)

$$D_{ijk} = p_{ijk} - p_i p_j p_k - p_i R_{jk} - p_j R_{ik} - p_k R_{ij};$$
 (12)

X the positive root of the equation

$$C_{ijk}(X^2 - 1) - D_{ijk}X = 0;$$
 (13)

$$F^* = 1/(X^2 + 1).$$
 (14)

However, we have K>3 fields to consider. There are $\binom{K}{3}$ choices of 3 fields each yielding X in (13). Accordingly, as a composite value for X we use the positive root of

$$(\sum_{ijk})(x^2-1) - (\sum_{ijk})x = 0,$$
 (15)

with summation over the $\binom{K}{3}$ choices, instead of (13). This step yields F^* . Notice that F^* (i.e.,X) is expressible in terms of the ratio D/C in (13), and the ratio $(\sum_{ijk})/(\sum_{ijk})$ in (15). We also define F_h^* based on (15) with summation restricted to

We also define F_h^* based on (15) with summation restricted to triplets (i,j,k) for which h=i, j or k. In other words, F_i^* is a composite F corresponding to field i specifically. In our example we have $F^*=.0503$, and values of F_i^* equal to

With F_0 = only .00338 we regard these values as evidence of positive field dependence.

Given the value of F, under independence we may show

$$R_{ij} = F(1-F)x_ix_j . \qquad (17)$$

with

$$y_i = p_i + (1-F)x_i \text{ and } u_i = p_i - Fx_i.$$
 (18)

Hence we now want to fit the values x_i and x_j . In (18) there are $\binom{K}{2}$ equations in K unknowns; for K=3 Fellegi and Sunter in effect solve the system of 3 equations in 3 unknowns. For K>3, from (17) we may form the following composite equation:

$$\log x_i = Z_i/(K-2) - Z/(K-1)(K-2) - .5 \log F_i^*(1-F_i^*)$$
 (19)

with Z_i the sum of log R_{jk} over the K-1 pairs (j,k) for which j = i or k = i, and Z the sum of log R_{jk} over all $\binom{K}{2}$ pairs, and with F** substituted for F specifically for field i.

Thus we use composite x_i (>0) based on (19). Then we get, from (18),

$$y_i^* = p_i + (1 - F_i^*) x_i$$
 and $u_i^* = p_i - F_i^* x_i$. (20)

These y_i^* and u_i^* , like F_i^* , are in essence just averages of Fellegi-Sunter 3-field results for the $\binom{K-1}{2}$ triplets which include i. In our example we get

Along with F^* and F_1^* too large, we obtain y_1^* too small (it should be much closer to 1) under the assumption of independence. (Yet it would be especially difficult to get the true value of y_1^* by an external sample of the true matches, because with true $F \leq F_0$, i.e., F less than .00338, these true matches are so scanty.)

We have based the development of (20) on the largest aggregates of agreement fractions: p_i , p_{ij} , p_{ijk} . This fact is important because for many fields even p_i is much closer to 0 than to 1, as our example indicated in (7); an aggregate such as p_{ijkmnh} may be so close to 0 as to be quite uninformative. In our example, Table 1 shows that p_{12345} is only .001266.

Thus we have fitted results for K fields under the assumption of independence. We now adjust y_i^* and u_i^* in preparation for the

next section: there we will be using the fraction \mathbf{F}_0 , as a device which permits us to fit interaction terms that reflect positive dependence. Let \mathbf{y}_i^* and \mathbf{u}_i^* be given by the 2x2 system of equations

$$y_{i}^{\dagger}(1-u_{i}^{\dagger})/u_{i}^{\dagger}(1-y_{i}^{\dagger}) = y_{i}^{*}(1-u_{i}^{*})/u_{i}^{*}(1-y_{i}^{*})$$
 (22)

$$F_0 y_i' + (1 - F_0) u_i' = p_i$$
 (23)

The requirement (22) preserves, up to a constant factor, the values of all the likelihood ratios U_S : that is, for two different cells S and T the ratio U_S/U_T is unchanged. (The hierarchical relationship among likelihood ratios is thus maintained.) We thus preserve the values of the weights $\mathbf{w_i}$ that appear in the constant-plus-sum relationship of (4) (although the value of the constant is altered). In this sense we maintain the spacing between y and u that was obtained from (16): that is, in this sense we hold effect (b) of Sec. 3 constant. We also ensure $0 < \mathbf{u_i'} < \mathbf{y_i'} < 1$. The requirement (23) ensures that, in Sec. 4, the fitted overall proportion of agreement for field i equals the observed $\mathbf{p_i}$.

One easily sets up and solves a quadratic equation to obtain y_i^* and u_i^* based on (22) and (23). In our example we get

4. Interactions. This section fits parameters to reflect positive dependence among field agreements, while preserving the values of \mathbf{U}_S obtained in Sec. 3. Consider the fitted probability

$$P_{S}' = F_{0}[_{i}\pi_{S}y_{i}'][_{i}\pi_{S}(1-y_{i}')] + (1-F_{0})[_{i}\pi_{S}u_{i}'][_{i}\pi_{S}(1-u_{i}')].$$
 (25)

With P_S^{τ} as starting point, consider G_S and P_S^{τ} given by

$$\log G_{S} = \log P_{S}^{\prime +} \sum_{i \leq j} b_{Sij} B_{ij}$$
 (26)

$$P_S^{\pi} = G_S/G \text{ with } G = \sum G_S.$$
 (27)

For $P_S' = 0$ we set $P_S'' = 0$. Here, P_S'' is an update of P_S' (equal to P_S' if all B_{ij} are 0). The quantities B_{ij} are interaction parameters, to be fitted to each cell as a whole. (Here we are thinking of y_i' and u_i' as fitted "main effects," which in this section are fixed.) For i<j, the constant b_{Sij} equals 1 for either (a) is and js or (b) is and js; it equals -1 for either (a) is and js or (b) is and js.

(Along with P" we also are fitting Q" and Q" according to (26) and (27), based on replacement of P', in (26) by Q's and Q's respectively. The scaling in (27) forces Q"s and Q"s, like P", to sum to 1 over S. Because we fit interaction parameters for cells as wholes rather than for matched and unmatched pairs separately, the values of all ratios U_S are preserved.)

Values $B_{ij}=0$ correspond to independence of events E_i and E_j , values $B_{ij}>(<)$ 0 to positive (negative) dependence. To correspond to the typical pattern of nonnegative dependence, we impose the restrictions $B_{ij} \ge 0$. Subject to these restrictions, we fit B_{ij} so as to maximize the likelihood function $\sum P_S \log P_S^n$, i.e., to

minimize
$$f = -\sum_{S} P_{S} \log P_{S}^{"}$$
 (28)

as a function of the variables ${\tt B}_{i\,j}$. As a measure of closeness of fit of (the minimizing) ${\tt P}_S^n$ to the original ${\tt P}_S$, we use

$$f^* = -\sum_{S} \log P_S^n + \sum_{S} \log P_S, \qquad (29)$$

which is always nonnegative. (Here 0 log 0 is 0.) We would obtain $f^* = 0$ only for a perfect fit: $p_S^* = P_S$ for all S. (We thus use the maximum likelihood criterion to measure closeness of fitted table proportions to given proportions. We could just as well use minimum Pearson chi-square, with results not greatly

different, with somewhat more complexity in the required computations.)

It is straightforward to show (Fiacco and McCormick 1968) that in (28) we have a convex programming problem. Suppose that we find a set of values B_{ij} (≥ 0) such that we have for i<j, always, either (a) $df/dB_{ij} = 0$ or (b) $B_{ij} = 0$ and $df/dB_{ij} > 0$. Then for this set of values, we have obtained a solution to the minimization problem in (28); and this solution is unique. A stopping rule may be based on these conditions (a) and (b); other computation issues are discussed in Sec. 8.

For us, a unique solution will routinely exist if the number of cells with $P_S>0$ comfortably exceeds the number of parameters (in all, 2K + K(K-1)/2) to be fitted. This requirement needs to be met, furthermore, for proper identifiability of parameters. We will not single out cases where there is no solution; but we would at least require K=5. Sec. 7 further discusses the choice of (fields and a value for) K.

In our example the difference f^* in (29) was .007143. Positive interactions were:

$$B_{23} = .0073$$
, $B_{34} = .1028$, $B_{15} = .0647$, $B_{25} = .0004$, $B_{45} = .0270$. (30)

With B_{ij} applied to each cell as a whole in (26) and (27) we can preserve the value of U_S and w_i , and thus maintain the relationship in (4) and the hierarchical relationship discussed in Sec. 1. We use only y_i^* and u_i^* as determined in Sec. 3 to determine the likelihood ratio U_S mainly because, in trying to minimize f, we would no longer have a convex problem with its corresponding unique global minimum, if we tried to take into account variation in U_S . Accordingly, we use the scheme of Sec. 3 to fit U_S .

However, we want to be able to apply this scheme of Sec. 3 realistically, under the assumption of mutual independence. Thus having fitted B_{ij} , our next step is, in effect, to remove interaction effects from our original table frequencies P_S - before going back to Sec. 3. Our new set of frequencies will

provide an analogue to a deseasonalized time series from which one has removed seasonal effects. For values $B_{ij} \ge 0$ obtained from the solution to the convex problem, we consider G_S and P_S^* given by

$$\log G_S = \log P_S - \sum_{i \le j} b_{Sij} B_{ij}$$
 (31)

$$P_S^* = G_S/G \text{ with } G = \sum G_S,$$
 (32)

in analogy with (26) and (27). For $P_S=0$ we set $P_S^*=0$. Then, P_S^* is our de-interactionized set of cell frequencies and hopefully tends to fit the model for P_S^* in (25) (without any interaction terms).

Recomputation. This section considers iterative alternation between the procedures in Sec. 3 and Sec. 4. The just-obtained table P* might behave as though (with F = F_0) the assumption of mutual independence in fact held. Using P* (rather than P_S), we go back and recompute y' and u' in (22), at the end of Sec. 3. We also recompute F* as in (15); if the new F* (while > F_0) is not less than the previous F*, we would regard the latest attempt to fit interaction terms, and remove additional positive dependence, as unsuccessful. We also view it as unsuccessful if we should obtain y* or u* outside the range <0,1>, or R_{ij} \leq 0 (in which case we can do little but stop).

If we are successful thus far, we fit a new set of interaction terms as follows. From the new y_i' and u_i' we form a new P_S' as in (25), except that to provide correspondence to P_S we add back the already obtained interaction terms — in the manner of (26) and (27), where P_S'' was formed. We then have a new starting point P_S' , from which to try to remove additional positive dependence. In other words, we are now at (25), at the start of Sec. 4.

New, additional interaction terms, and new $P_S^{"}$, are then fitted using (26) and (27). We get a new minimal value of f as in (28), and a difference as in (29), measuring discrepancy between $P_S^{"}$ and P_S . If the value of (29) is greater than the

previous value, however, we regard the whole attempt to refit y_1' , u_1' and P_S^m as unsuccessful. At this point, if we have been unsuccessful in any sense, we go back to the <u>previous</u> set of y_1' and u_1' , and use the accompanying w_1 and P_S^m as final values.

However, if we are successful, we give it another try. Based on the accumulated sums of interaction terms fitted thus far, we de-interactionize, re-fit P_S' , re-compute f (and an additional set of interaction terms), etc. We can keep going until we are unsuccessful, and then revert to results for the previous fit. Thus we consider results based on the best available fit for our 2nd-order model: y_i' , u_i' and cumulative $B_{i,j}$ based on them.

(Quite possibly, straightforward variants of some of the above-discussed computational steps will be helpful for particular sets of data, in getting f as small as possible, subject to the model in (26) and (27). The exact computational steps are thus somewhat open to investigation. However, we have found that some of the more obvious variants do not seem helpful. In any event we would preserve the essential, alternating calculations: in Sec. 3 extend the Fellegi-Sunter method of moments, and in Sec. 4 fit second-order nonnegative interaction terms based on a convex programming problem, so as to maintain the likelihood-ratio relationship in (4).)

In our example we were successful until on the 14th set of computations for y_i^* and u_i^* we obtained an $R_{i,j} \le 0$. Hence we consider the results from the 13th set of calculations. We got 13th $F^* = .0147$, vs. initial $F^* = .0503$, and vs. $F_0 = .00338$; hence most of the discrepancy from F_0 has been removed, and most of the positive dependence between fields may have been removed. Values of F_i^* also decreased similarly. The value of the difference f^* in (29) was reduced to only .002440, vs. .007143 after the 1st cycle; hence in a sense our iterations have removed most of the initial discrepancy between observed and fitted cell proportions. Accumulated interactions after the 13th cycle were (cf. initial values in (30))

$$B_{12} = .0351$$
, $B_{23} = .1111$, $B_{14} = .0644$, $B_{34} = .2556$, (33) $B_{15} = .1172$, $B_{25} = .0082$, $B_{45} = .2589$.

For the 13th cycle we also obtain (compare to initial values in (24) which correspond to the assumption of independence)

In (34), field 1 (especially) and fields 2 and 5 are de-emphasized, while fields 3 and (especially) 4 are given more emphasis, than in (24). Moreover, we now have $w_4 > w_3$ in spite of the facts (a) $p_3 < p_4$ in (8), indicating that field 3 might be a more useful discriminator than field 4, and (b) initially $w_3 > w_4$.

We would use these results for the data of Table 1. In this example a better fit was prevented only by $R_{i\,j}$ becoming nonpositive.

Iteration in this example was beneficial, with 13 productive cycles. In other examples there will be fewer (or more) such cycles, perhaps only 1; but even a single productive cycle with fitted interactions is an advance from the independence model of Sec. 3.

6. Final Results. At this point we have a clear choice for w_i and U_S : the last set of values for these, for which success was achieved. However, we want also to fit the best legitimately constructed final values for Q_{1S} and Q_{2S} . In this section we obtain these values, and then consider posterior probabilities based on them: for each cell S, a conditional probability that (a pair belonging to) the cell represents a true match.

We consider a 2-dimensional table. There will be 2 rows, corresponding to match and non-match, and 2^K columns, corresponding to the 2^K cells S. As in Sec. 4 we form $Q_{1S}^{"}$ in the manner of $P_{S}^{"}$ except that instead of log $P_{S}^{'}$ in (26) we use log $Q_{1S}^{'}$; we form $Q_{2S}^{"}$ similarly. Then we put $F_{0}Q_{1S}^{"}$ in row 1, column S and $(1-F_{0})Q_{2S}^{"}$ in row 2, column S. But we want, along

with row sums equal to F_0 and $1-F_0$, to have column sums equal to the observed P_S : having completed the fit of interactions, we now make full use of the observed proportions, P_S . We use iterative proportional fitting (Ireland and Kullback 1968) or "raking" to produce a new table which is minimally distorted from the original $2x2^K$ table, subject to: (a) row and column sums having the above totals; and (b) in the manner of (22) toward the end of Sec. 3, preservation of the constant-plus-sum relationship in (4) and the values of w_i in this relationship. When the raking is completed, we have table entries of form (row 1) $F_0Q_{1S}^*$ and (row 2) $(1-F_0)Q_{2S}^*$ with Q_{1S}^* and Q_{2S}^* , each summing to over S, representing final fitted probabilities.

Let M denote the event "match" and $U_S^* = Q_{1S}^*/Q_{2S}^*$. Given that prior P(M) is F_0 , and P(S|M) is Q_{1S}^* , with P(S) = P_S , we have the posterior probability

 $P(M|S) = F_0Q_{1S}^*/P_S = F_0U_S^*/(F_0U_S^* + 1 - F_0). \tag{35}$ (Note that P_S equals $F_0Q_{1S}^* + (1 - F_0)Q_{2S}^*$, in this development.) Thus we merely divide the fitted cell entry in row 1 and column S, by the column total P_S , to obtain P(M|S). As indicated in Sec. 1, we would make use of these posterior probabilities P(M|S). Note that if P_S is 0 we do not have a meaningful P(M|S), but that an empty cell does not matter anyway. Table 2 gives values of P(M|S), in the fashion of Table 1, for our example. Wherever there is agreement for fields 3 and 4, we have P(M|S) > .98; otherwise it is <.5. Using a cutoff value of .90 or .95, we would for this example be inclined to view agreement in these two fields as evidence of true match, and disagreement in either of them as lack of evidence of it.

	Tal	able 2. Values of		P(MS)	
00000 .00005	0 00010	.096580	00001	.000119	00011 .202047
10000 .00005	4 10010	.103233	10001	.000128	10011 .214242
01000 .00013	6 01010	.224579	01001	.000323	01011 .406871
11000 .00014	7 11010	.237729	11001	.000347	11011 .424847
00100 .02397	5 00110	empty	00101	.054982	00111 .991980
10100 .02577	0 10110	.982527	10101	.058957	10111 .992547
01100 .06239	6 01110	.992981	01101	.136160	01111 .997024
11100 .06686	9 11110	.993478	11101	.145101	11111 .997236

The purpose of this paper, however, is more than to reach such a conclusion for our example. As we indicated in Sec. 1, our chief goal is to provide the best possible values for Q_{1S} and Q_{2S} . These values may then be put to use in a hypothesis-test framework, for the cost-utility approach of Tepping (1968), or in other ways. Or one could solve a transportation problem based on P(M|S) as in Sec. 1.

As indicated in Sec. 1, the value of F_0 in (2) represents an upper bound on the true value of F rather than the true value itself; we would carry out our calculations for smaller values of F, to see how the results change. We have done this for P(M|S); but one can, of course, do similar investigations for the hypothesis-test framework. One can surmise from (35), that use of a smaller value for F_0 and thus a smaller prior P(M) should yield smaller values for P(M|S); and in our investigation this was the case.

For this investigation we replaced F_0 in (2) by p_{34} , the proportion of all pairs with agreement in fields 3 and 4, in view of the Table 2 results which suggested the decision "match" if and only if there is agreement for both these fields. The value of p_{34} is .001731, based on numerator 2287 (vs. 4465 for F_0). In the sense of Sec. 5 we were unsuccessful (F^* failed to decrease) on the 7th cycle, so we used 6th-cycle results as given in Table 3. There is still a marked dichotomy between cells S for which fields 3 and 4 agree, and all other cells; but Table 3 provides a reminder that P(M|S) in Table 2 is overstated and that agreement in fields 3 and 4 does not guarantee that we have a true match.

016 0001	0.001170				
	0 .003472	00001	.000068	00011	.014637
1001	0 .004787	10001	.000094	10011	.020095
051 0101	0 .011011	01001	.000218	01011	.045314
71 1101	0 .015137	11001	.000301	11011	.061495
115 0011	0 empty	00101	.013145	00111	.743760
295 1011	0 .484505	10101	.018056	10111	.800279
384 0111	0 .685084	01101	.040823	01111	.902673
594 1111	0 .750199	11101	.055494	11111	.927555
3	84 0111	84 01110 .685084 94 11110 .750199	84 01110 .685084 01101 94 11110 .750199 11101	84 01110 .685084 01101 .040823 94 11110 .750199 11101 .055494	84 01110 .685084 01101 .040823 01111 94 11110 .750199 11101 .055494 11111

One might wonder how to choose a "best" value for F that is $\leq F_0$. One could consider the final value of F* in (29): choose the F that minimizes this objective function. But because our model does not fully permit us to account for all positive dependence (source (c) of Sec. 2), it seems that we will fit a value of F which tends to be too large (source (a) of Sec. 2) and are apt to overstate the value of F. Moreover, just as for fitting best values for ratios U_S , we encounter difficulty because we no longer have a convex programming problem. (Whereas for F = F_0 we obtained, as in Sec. 5, a final f^* = .002440, we get f^* = .004220 for F = P_3 4. That is, F_0 corresponds to a better fit than does P_3 4.)

A final note: with E_i denoting the event "agreement for field i," we may sum over Q_{1S}^* for isS, to get final fitted values of $P(E_i \mid M)$ for i=1, ..., K. In our example these are

These are not as close to 1 as they realistically ought to be at least partially because the maximal F_0 exceeds the true fraction of matches. When we replace F_0 by p_{34} as above the fitted values of $P(E_i \mid M)$ increase to

7. Choice of Fields. In this section we consider how best to form fields. From Sec. 1, recall that field 1 in our example is based on 2 data items (race and ethnicity). Likewise fields 2, 3 and 4 are each based on 2 items, while field 5 is based on 3 items. In all, we have 11 "items" although we have been only using 5 fields. One might ask, why not form and use 11 fields? (Moreover, by using binary digits for each item, we could form many more than 11 fields.)

Accordingly, we now consider the general issue of how best to form fields from a given set of data items. As a first principle, we would include every item in one (and only one) data

field, in order to make full use of available information concerning agreement/disagreement patterns. But for the following reasons we would not necessarily form a field from each single item. (That is, we would not necessarily form 11 fields in our example.)

- (a) As the number of fields (K) increases, there is increasing room for complicated interactions among the fields. Departures from our second-order model in Sec. 4 may be exacerbated.
- (b) By increasing K, we do increase the number of fitted parameters w_i , and B_{ij} . However, we do not necessarily get more information. For example, if in our illustration we split raceethnicity into two separate fields, we get two separate weights w_i , and a suggestion of more information. However, we do not get a single weight for the combination race-ethnicity, that reflects the joint behavior of the two items. Nor do we get a pairwise interaction, B_{ij} , which corresponds to this combination, as a single field, in conjunction with a second field such as birthdate.
- (c) (less vital) As in the next section, the required order of computations blows up as K increases.

We would prefer K=5, provided that $P_S>0$ for all or nearly all of the 32 cells S, in order to permit comfortable fitting of 20 parameters as in Sec. 4. There are, however, two further possible objections to small K. We state and deal with these as follows.

(a) For K (and 2^K , the number of cells) small, there can be awkwardness, because of discreteness, in constructing a decision rule analogous to the classical hypothesis test (Sec. 1). Cells are few in number with large probabilities Q_{1S} and Q_{2S} associated with them; as a result one typically cannot come close to attaining the desired Type I and Type II error probabilities. But as Fellegi and Sunter indicate, one can introduce an artifical randomizing device to attain these. Moreover, in Sec. 1 we indicated how to use posterior probabilities either to solve

a transportation problem or to construct decision rules based on cutoffs instead of the hypothesis-test framework. Hence this objection can be overcome.

(b) The proportions of agreement, p_i (as in (8)), are smallest for the smallest K and the most aggregation of items. Small values of p_i can lead to difficulties when the denominator ("D") of p_i is small: that is, when $\sum N_{hA}N_{hB}$ in Sec. 1 is small. For this reason one might thus anticipate difficulty for small K. However, the rest of this section indicates that such difficulty can be avoided for K as small as 5.

For D = 10 million and p_i = .0002 we have a numerator 2000. If numerators of p_i are, in this fashion, in the thousands, then we have relatively small sampling error and may usefully compare p_i to each other - and likewise utilize p_{ij} and p_{ijk} as we have done for our example. On the other hand, suppose D is 10,000; for p_i = .0002 we are dealing with a numerator 2. Sampling error, extreme discreteness, etc., creates an intractable situation. For sufficiently large D, with numerators large enough, there should be no objection to having a small K and small values of p_i . (For given files A and B, the numerators will increase as D increases, although they will not proportionately keep pace - because the overall number of true matches, which relatively account for more agreements than nonmatches, stays the same while the overall number of nonmatches increases. In any event, it seems best to choose fields so that the p_i 's are equalized as nearly as possible, as a "maximin" strategy.)

Thus we next must consider how big a D we want to deal with, and thus how many blocks we want to form: the fewer the blocks, the bigger D will be, for given files A and B. For K fields we need to make DK field comparisons as to field agreement/disagreement, in order to obtain the quantities P_S . These comparisons can be streamlined, and based on just integers. Thus computationally it becomes quite feasible to consider D equal to, say, 10 million (at least) in conjunction with K=5, in striving to make the numerators for p_1 , p_{11k} acceptably large.

Another reason for forming a few large blocks, as opposed to many small blocks, is that likelihood of error is reduced, in the following sense. Recall from Sec. 1 that blocking is to be done so that: if a pair of records from A and B constitute a true match, they belong to the same block. Blocking leads to "error" if it assigns two such records to different blocks. By having blocks as few in number and as large as possible, we reduce the possibilities for such error.

8. Computation. This section pertains strictly to computational issues for the convex programming problem of Sec.
4. For this general problem, of course, numerous algorithms have been developed; but we would like to add the following.

We start each (convex programming) problem with B_{ij}, corresponding to additional positive interaction, set to 0. To compute Vf, the vector of 1st derivatives, and $\nabla^2 f$, the matrix of 2nd derivatives, for f in (28) with respect to B_{ij}, we need, in the notation for (26), the sums $\sum_S b_{Sij}G_S$. We also need, for $i \neq j \neq k \neq i$, analogous sums $\sum_S c_{Sijk}G_S$ which correspond to 3rd-order interaction; likewise we need sums which correspond to 4th-order interaction. In all, the number of required additions involving G_S is seen to be of order

$$2^{K-1} \int_{1-2}^{4} {K \choose j},$$
 (37)

which increases dramatically as K increases: important because of the need to compute these derivatives repeatedly. Thus, as indicated in Sec. 7, there can be computational problems for large K.

A final computational aid for persons familiar with convex programming algorithms: suppose that the direction vector,

$$(\nabla^2 r)^{-1}(\nabla r), \tag{38}$$

has a positive component corresponding to a B_{ij} which is 0. Minimizing f along this direction vector, we would be led to make B_{ij} negative, contrary to the constraints B_{ij} ≥ 0 . Under these

conditions we have recomputed the direction vector with such ${\bf B_{ij}}$ excluded (temporarily) from the set of variables. If necessary, we repeat to make further exclusions. This approach permits us neatly to bound ${\bf B_{ij}} \ge 0$ without the use of penalty functions; and we find that we quickly get to the solution. For a similar reason, when there are positive components for ${\bf B_{ij}} > 0$, we multiply by a positive scalar small enough to keep all ${\bf B_{ij}} \ge 0$.

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