# Poisson approximation in connection with clustering of random points

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#### Abstract

Let n particles be independently and uniformly distributed in a rectangle  $\mathbf{A} \subset \mathbb{R}^2$ . Each subset consisting of  $k \leq n$  particles may possibly aggregate in such a way that it is covered by some translate of a given convex set  $C \subset \mathbf{A}$ . The number of k-subsets which actually are covered by translates of C is denoted by W. The positions of such subsets constitute a point process on  $\mathbf{A}$ . Each point of this process can be marked with the smallest necessary "size" of a set, of the same shape and orientation as C, which covers the particles determining the point. This results in a marked point process.

The purpose of this paper is to consider Poisson (process) approximations of W and of the above point processes, by means of Stein's method. To this end the exact probability for k specific particles to be covered by some translate of C is given.

#### 1 Introduction

Assume n particles are uniformly and independently distributed in a rectangle  $\mathbf{A} \subset \mathbb{R}^2$ , and let  $C \subset \mathbf{A}$  be a convex set, small relative to  $\mathbf{A}$ . To avoid problems with the boundaries of  $\mathbf{A}$ , the torus convention will be used throughout the paper. The aim of this paper is to investigate Poisson approximation of certain variables and point processes which concern subsets of particles which are covered by some translate of C.

First a brief review of related problems will be given. Thereafter necessary notation and some general Poisson approximation theorems are presented. In Section 3 the probability that a fixed number of particles aggregate in such a way that they are all covered by some translate of C is given, and its connection with integral geometry is discussed.

Section 4 constitutes the bulk of the paper. First the number of subsets consisting of k < n particles, called k-subsets in the following, which are covered by some translate of C is considered. The Stein-Chen method is used to bound the error made in the approximation of this number by a Poisson variable.

To each k-subset, which is covered by some translate of C, can be attached its position on A, and the smallest  $s \in \mathbb{R}^+$  for which some translate of sC covers the k particles. Poisson process approximation of three point processes determined by these positions and sizes is also dealt with in Section 4, as is the joint distribution of the number of covered k-subsets, and the corresponding processes, for different k. The next section deals with asymptotic aspects, such as rates of convergence. The paper is concluded with some problems of possible interest for more detailed investigation.

Note that the parameters used in the approximations follow from results in Section 3. For those readers whose main interest is the approximations, and not the geometrical considerations leading to the parameters, it is enough to use Section 3 as a reference.

#### 1.1 A brief historical account

There are  $\binom{n}{k}$  different k-subsets of particles, some of which are covered by translates of C. Order the k-subsets in some way and let W denote the number of k-subsets which actually are covered

by some translate of C. Then W can be written as

$$W = \sum_{i=1}^{\binom{n}{k}} I_i,$$

where

$$I_i = \begin{cases} 1 & \text{if the } i \text{th } k \text{-subset is covered by some translate of } C, \\ 0 & \text{otherwise.} \end{cases}$$

A k-subset which is covered by some translate of C will in this introduction be referred to as a k-aggregate in accordance with previous literature. Silberstein (1945), Mack (1948,1949), and Eggleton and Kormack (1944) are concerned mainly with the calculation of the expected number of k-aggregates, E[W], in one dimension (then C obviously is a line segment) and, for certain shapes of C, in two dimensions. Circular sets are handled by Mack (1948) and Silberstein (1945), while rectangular sets are handled by Mack (1949) and Eggleton and Kormack (1944). Mack (1949) discusses arbitrary shapes in two dimensions, and also gives some results in higher dimensions, and Mack (1948) argues for a Poisson limit of the number of k-aggregates as n tends to infinity.

The cases of a disc and a square are treated by Aldous (1989). He gives approximate formulas for the distribution of the radius (side length) of the smallest disc (square) which can be translated to a position where it covers at least k particles.

In the special case of k=2 and a circular set, C, the number of k-aggregates equals the number of pairs of particles with interpoint distance less than the diameter of C. Convergence of this number to a Poisson limit is discussed in Silverman and Brown (1978). The results therein are complemented with a bound for the total variation distance in Silverman and Brown (1979). In e.g. Kryscio and Saunders (1983) Poisson convergence of the interpoint distances is considered in the case where the underlying process of particles is not necessarily stationary.

In Silverman and Brown (1978) the number of close pairs is given as an example of the classical U-statistics. Poisson approximation of U-statistics and sums of dissociated variables (defined in Section 2.2) are treated by Barbour and Eagleson (1984) and Barbour, Holst and Janson (1992). In these two references the Stein-Chen method is used to bound the total variation distance for Poisson approximations of the sums; in particular, the bound in Silverman and Brown (1979) is improved.

The number of close pairs is sometimes used to test the null hypothesis that a sample is uniformly distributed against some clustering or regular alternative. This is discussed e.g. in Silverman and Brown (1978) and in Barbour, Holst, and Janson (1992) p. 34.

Another related concept is the so-called *scan statistic*, i.e. the maximal number of particles which are covered by some translate of C. In this context the particles usually constitute a Poisson process rather than being a fixed number. In one dimension considerable attention has been devoted towards finding approximations of the distribution of the scan statistic, see e.g. Naus (1982), Alm (1983), Janson (1984), and Glaz (1989). The references on scan statistics in higher dimensions are more scattered. The two-dimensional case is treated in Alm (1997) and Loader (1991), the latter, however, being restricted to rectangular sets.

In one dimension the scan statistic is used in the same way as the number of close pairs is used in two dimensions, in testing the null hypothesis of an underlying Poisson process against a clustering or regular alternative.

#### 2 Preliminaries

#### 2.1 Notions and notation

Let  $\mathbb{R}^2$  denote the 2-dimensional Euclidean space, with a fixed origin, O, and orthogonal coordinate-axes. The *area* of a (measurable) subset of  $\mathbb{R}^2$  is its 2-dimensional Lebesgue measure which we denote by  $\mu$ .

For  $B, C \subset \mathbb{R}^2$  and  $c \in \mathbb{R}$  the Minkowski sum and scalar multiple are defined as

$$B + C = \{x + y : x \in B, y \in C\}$$
 and  $cB = \{cx : x \in B\},\$ 

respectively. If c = -1 we get  $\check{B} = \{-x : x \in B\}$ , which we call the reflected set of B. For  $x \in \mathbb{R}^2$ ,  $B + \{x\}$  is the translate of B by x, which is denoted by B + x. If  $B = \check{B} + x$  for some  $x \in \mathbb{R}^2$ , B is said to be centrally symmetric. An alternative, and for us more useful, way of writing the Minkowski sum is

$$B + C = \{x : B \cap (\check{C} + x) \neq \emptyset\}. \tag{2.1}$$

For the set xB + yC, where  $x, y \in \mathbb{R}^+$  and  $B, C \subset \mathbb{R}^2$  are non-empty convex sets, the area can be written as:

$$\mu(xB + yC) = x^2\mu(B) + 2xy\nu(B, C) + y^2\mu(C), \tag{2.2}$$

where  $\nu(B,C)$  is the *mixed area* of B and C, which is actually defined by (2.2) (see e.g. Bonnesen and Fenchel (1948) p. 40). These mixed areas are usually only defined for non-empty sets. We make the natural extension that if B or  $C = \emptyset$ , then  $\nu(B,C) = 0$ . It can be shown that if C is a convex set in  $\mathbb{R}^2$  then

$$\mu(C) < \nu(C, \check{C}) < 2\mu(C), \tag{2.3}$$

where the lower bound is attained if and only if C is centrally symmetric while the upper bound is attained if and only if C is a triangle. For a proof of these facts, see e.g. Bonnesen and Fenchel (1948) p. 105.

Assume that the rectangle in which the particles are distributed,  $\mathbf{A}$ , is centred at the origin. The family of convex sets,  $C \subset \mathbb{R}^2$ , with the properties that  $C + \check{C} \subset \mathbf{A}$  and that O is an interior point of C, we denote by  $\mathcal{K}$ .

Let  $(\mathcal{X}, \mathcal{A})$  be any measurable space. The total variation distance,  $d_{TV}$ , between two probability measures  $\mu$  and  $\nu$  on  $\mathcal{X}$  is defined to be

$$d_{TV}(\mu,\nu) = \sup_{A \in A} |\mu(A) - \nu(A)|.$$

If the state space is discrete, then

$$d_{TV}(\mu, \nu) = \frac{1}{2} \sum_{i \in \mathcal{X}} |\mu\{i\} - \nu\{i\}|,$$

and in this case convergence in total variation distance,  $d_{TV}(\mathcal{L}(X_n), \mathcal{L}(X)) \to 0$ , is equivalent to  $\{X_n\}$  converging in distribution to X.

#### 2.2 General Poisson approximation theorems

Following the notation of Barbour, Holst and Janson (1992), let  $\Gamma$  be an arbitrary finite collection of indices and let

$$W = \sum_{\alpha \in \Gamma} I_{\alpha}, \quad \pi_{\alpha} = E[I_{\alpha}] \quad \text{and} \quad \lambda = E[W],$$

where  $I_{\alpha}$ ,  $\alpha \in \Gamma$ , are, possible dependent, indicator variables. The following method for bounding the total variation distance between the distribution of  $\mathcal{L}(W)$  and a Poisson variable with parameter  $\lambda$  is called the local approach of the Stein-Chen method. It is suitable to use when there is a natural dependence structure which allows every pair of indicators to be classified as either strongly or weakly dependent. For each  $\alpha \in \Gamma$ , let  $\Gamma \setminus \{\alpha\}$  be divided into two subsets; one consisting of those  $\beta \in \Gamma$  for which  $I_{\beta}$  is weakly dependent on  $I_{\alpha}$ , and one consisting of the indicators which are strongly dependent on  $I_{\alpha}$ . Denote these subsets by  $\Gamma_{\alpha}^{w}$  and  $\Gamma_{\alpha}^{s}$ , respectively, and let

$$Z_{\alpha} = \sum_{\beta \in \Gamma_{\alpha}^{s}} I_{\beta}.$$

**Theorem 2.1** (Theorem 1.A Barbour et al. (1992)). Let  $\Gamma$  be an arbitrary finite collection of indices. With the above definitions, for any choice of the index sets  $\Gamma^s_{\alpha}$ ,  $\alpha \in \Gamma$ ,

$$d(\mathcal{L}(W), \operatorname{Po}(\lambda)) \leq \sum_{\alpha \in \Gamma} (\pi_{\alpha}^{2} + \pi_{\alpha} E[Z_{\alpha}] + E[I_{\alpha} Z_{\alpha}]) \lambda^{-1} (1 - e^{-\lambda})$$
$$+ \sum_{\alpha \in \Gamma} \eta_{\alpha} \min(1, \lambda^{-1/2}),$$

where

$$\eta_{\alpha} = E[|E[I_{\alpha} \mid (I_{\beta}, \beta \in \Gamma_{\alpha}^{w})] - \pi_{\alpha}|].$$

The next theorem concerns approximation of marked point processes. Let  $\Xi = \sum_{\alpha \in \Gamma} I_{\alpha} \delta_{Y_{\alpha}}$ , where  $\delta_y$  denotes the unit point mass at y and the  $Y_{\alpha}$ 's are random variables with state space  $\mathcal{Y}$ , which is assumed to be metric and separable. Then  $\Xi$  is a random element on the space  $\mathcal{Z}$  of configurations of finite point processes over  $\mathcal{Y}$ . Assume that a measure on  $\mathcal{Y}$  is defined for each  $\alpha \in \Gamma$  by

$$\lambda_{\alpha}(A) = P(I_{\alpha} = 1, Y_{\alpha} \in A) \tag{2.4}$$

and let

$$\lambda = \sum_{\alpha \in \Gamma} \lambda_{\alpha}.$$

Furthermore, let

$$\Xi^w_lpha = \sum_{eta \in \Gamma^w_lpha} I_eta \delta_{Y_eta} \quad ext{ and } \quad \Xi^s_lpha = \sum_{eta \in \Gamma^s_lpha} I_eta \delta_{Y_eta}.$$

The following theorem is the "local" version of Theorem 10.E in Barbour et al. (1992). Its proof is a combination of the proofs of Theorem 10.A and Theorem 10.E of that reference, which should be consulted for more details.

**Theorem 2.2** Let  $\Xi$  and  $\lambda$  be defined as above. Then

$$d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda)) \leq \sum_{\alpha \in \Gamma} \left( \pi_{\alpha}^{2} + \pi_{\alpha} E[Z_{\alpha}] + E[I_{\alpha} Z_{\alpha}] + \int_{\mathcal{Y}} d_{TV}(\mathcal{L}(\Xi_{\alpha}^{w} \mid I_{\alpha} = 1, Y_{\alpha} = x), \mathcal{L}(\Xi_{\alpha}^{w})) \lambda_{\alpha}(dx) \right).$$

**Proof.** Let

$$(\mathcal{A}h)(\xi) = \int_{\mathcal{Y}} [h(\xi + \delta_x) - h(\xi)] \lambda(dx) + \int_{\mathcal{Y}} [h(\xi - \delta_x) - h(\xi)] \, \xi\{dx\},$$

 $\xi \in \mathcal{Z}$ , be the generator of the immigration-death process on  $\mathcal{Y}$  with immigration intensity  $\lambda = \sum_{\alpha \in \Gamma} \lambda_{\alpha}$  and unit per-capita death rate, whose equilibrium distribution is that of the Poisson process on  $\mathcal{Y}$  with intensity  $\lambda$ . Let  $h_{\lambda,B}$  be the solution of

$$(\mathcal{A}h)(\xi) = I(\xi \in B) - Po(\lambda)(B),$$

which can be found in Proposition 10.1.1 in Barbour et al. Then

$$d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda)) = \sup_{B \subset \mathcal{Y}} |E[(\mathcal{A}h_{\lambda,B})(\Xi)]|.$$

In order to bound  $|E[(Ah_{\lambda,B})(\Xi)]|$  we write

$$E[(\mathcal{A}h)(\Xi)] = E\left\{\sum_{\alpha\in\Gamma}\int_{\mathcal{Y}}\left[h(\Xi+\delta_{x})-h(\Xi)\right]\lambda_{\alpha}(dx) + \sum_{\alpha\in\Gamma}\int_{\mathcal{Y}}\left[h(\Xi-\delta_{x})-h(\Xi)\right](I_{\alpha}\delta_{Y_{\alpha}})\{dx\}\right\}$$

$$= \sum_{\alpha\in\Gamma}E\left\{\int_{\mathcal{Y}}\left[h(\Xi-\delta_{x})-h(\Xi)\right](I_{\alpha}\delta_{Y_{\alpha}})\{dx\}\right\} - E[I_{\alpha}\left(h(\Xi_{\alpha}^{w})-h(\Xi_{\alpha}^{w}+\delta_{Y_{\alpha}})\right)]\right\}$$

$$+ \sum_{\alpha\in\Gamma}\int_{\mathcal{Y}}E[h(\Xi+\delta_{x})-h(\Xi)]\lambda_{\alpha}(dx) + \sum_{\alpha\in\Gamma}\int_{\mathcal{Y}}E[h(\Xi_{\alpha}^{w})-h(\Xi_{\alpha}^{w}+\delta_{x})]\lambda_{\alpha}(dx)$$

$$+ \sum_{\alpha\in\Gamma}E[I_{\alpha}\left(h(\Xi_{\alpha}^{w})-h(\Xi_{\alpha}^{w}+\delta_{Y_{\alpha}})\right)] - \sum_{\alpha\in\Gamma}\int_{\mathcal{Y}}E[h(\Xi_{\alpha}^{w})-h(\Xi_{\alpha}^{w}+\delta_{x})]\lambda_{\alpha}(dx). \tag{2.5}$$

To bound the absolute value of the right-hand side of the above equation we need Lemma 10.1.3 in Barbour et al. which says that

$$\Delta_1 h = \sup_{\xi \in \mathcal{Z}: x \in \mathcal{Y}} |h(\xi + \delta_x) - h(\xi)| \le 1 \tag{2.6}$$

and

$$\Delta_2 h = \sup_{\xi \in \mathcal{Z}; x, y \in \mathcal{Y}} |h(\xi + \delta_x + \delta_y) - h(\xi + \delta_x) - h(\xi + \delta_y) + h(\xi)| \le 1, \tag{2.7}$$

where h is the solution of

$$(\mathcal{A}h)(\xi) = I(\xi \in B) - Po(\lambda)(B).$$

By rewriting the first term on the right-hand side of (2.5), noting that the number of points of  $\Xi_{\alpha}^{s}$  is  $Z_{\alpha}$ , and using (2.7) we get

$$\begin{split} &\left| \sum_{\alpha \in \Gamma} E[I_{\alpha} \left\{ h(\Xi - \delta_{Y_{\alpha}}) - h(\Xi) - h(\Xi_{\alpha}^{w}) + h(\Xi_{\alpha}^{w} + \delta_{Y_{\alpha}}) \right\}] \right| \\ &= \left| \sum_{\alpha \in \Gamma} E[I_{\alpha} \left\{ h(\Xi_{\alpha}^{w} + \Xi_{\alpha}^{s} + I_{\alpha} \delta_{Y_{\alpha}}) - h(\Xi_{\alpha}^{w} + \Xi_{\alpha}^{s} + I_{\alpha} \delta_{Y_{\alpha}} - \delta_{Y_{\alpha}}) - h(\Xi_{\alpha}^{w} + \delta_{Y_{\alpha}}) + h(\Xi_{\alpha}^{w}) \right\}] \right| \\ &\leq \left| \sum_{\alpha \in \Gamma} E[I_{\alpha} Z_{\alpha} \triangle_{2} h] \right| \\ &\leq \sum_{\alpha \in \Gamma} E[I_{\alpha} Z_{\alpha}]. \end{split}$$

Similarly the second term on the right-hand side of (2.5) is bounded by

$$\Big| \sum_{\alpha \in \Gamma} E[(I_{\alpha} + Z_{\alpha}) \triangle_{2} h] \int_{\mathcal{Y}} \lambda_{\alpha}(dx) \Big| \quad \leq \quad \sum_{\alpha \in \Gamma} E[I_{\alpha} + Z_{\alpha}] E[I_{\alpha}].$$

Finally, the third term on the right of (2.5) is equal to

$$\left| \sum_{\alpha \in \Gamma} \int_{\mathcal{Y}} \left\{ E[h(\Xi_{\alpha}^{w}) - h(\Xi_{\alpha}^{w} + \delta_{x}) \mid I_{\alpha} = 1, Y_{\alpha} = x] - E[h(\Xi_{\alpha}^{w}) - h(\Xi_{\alpha}^{w} + \delta_{x})] \right\} \lambda_{\alpha}(dx) \right|. \tag{2.8}$$

By (2.6)  $|h(\Xi_{\alpha}^{w}) - h(\Xi_{\alpha}^{w} + \delta_{x})| \leq 1$ , and hence, using the definition of total variation distance, (2.8) is bounded by

$$\sum_{\alpha \in \Gamma} \int_{\mathcal{Y}} d_{TV}(\mathcal{L}(\Xi^w_\alpha \mid I_\alpha = 1, Y_\alpha = x), \mathcal{L}(\Xi^w_\alpha)) \lambda_\alpha(dx),$$

and the theorem is proved.

If  $\Gamma$  is a collection of k-subsets of  $\{1, 2, \ldots, n\}$ , then the family  $\{X_{\alpha} : \alpha \in \Gamma\}$  of random variables is said to be *dissociated*, as defined by McGinley and Sibson (1975), if  $(X_{\alpha} : \alpha \in A)$  and  $(X_{\alpha} : \alpha \in B)$  are independent whenever  $(\bigcup_{\alpha \in A} \alpha) \cap (\bigcup_{\alpha \in B} \alpha) = \emptyset$ . In the following theorem families of indicator variables belonging to a somewhat wider group are considered; instead of just one value of k, we let  $\Gamma$  be a collection of arbitrary subsets of  $\{1, 2, \ldots, n\}$ .

**Theorem 2.3** (i) Let  $\Gamma$  be a collection of subsets of  $\{1, 2, ..., n\}$  and the family  $\{I_{\alpha} : \alpha \in \Gamma\}$  be such that  $(I_{\alpha} : \alpha \in A)$  and  $(I_{\alpha} : \alpha \in B)$  are independent whenever  $(\bigcup_{\alpha \in A} \alpha) \cap (\bigcup_{\alpha \in B} \alpha) = \emptyset$ . If  $\Gamma_{\alpha}^{s} = \{\beta \in \Gamma : \beta \neq \alpha, \beta \cap \alpha \neq \emptyset\}$ , then

$$d_{TV}(\mathcal{L}(W), \text{Po}(\lambda)) \leq \sum_{\alpha \in \Gamma} (\pi_{\alpha}^2 + \pi_{\alpha} E[Z_{\alpha}] + E[I_{\alpha} Z_{\alpha}]) \lambda^{-1} (1 - e^{-\lambda}).$$

(ii) Let  $\Gamma$  and  $\Gamma_{\alpha}^{s}$  be defined as above, and assume that the family  $\{I_{\alpha}\delta_{Y_{\alpha}}: \alpha \in \Gamma\}$  is such that  $(I_{\alpha}\delta_{Y_{\alpha}}: \alpha \in A)$  and  $(I_{\alpha}\delta_{Y_{\alpha}}: \alpha \in B)$  are independent whenever  $(\bigcup_{\alpha \in A}\alpha) \cap (\bigcup_{\alpha \in B}\alpha) = \emptyset$ . Then

$$d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda)) \leq \sum_{\alpha \in \Gamma} \pi_{\alpha}^2 + \pi_{\alpha} E[Z_{\alpha}] + E[I_{\alpha} Z_{\alpha}].$$

**Proof.** The result is an immediate consequence of Theorem 2.1 and 2.2, since by the definition of  $\Gamma_{\alpha}^{s}$  it follows that  $I_{\beta}$  and  $I_{\beta}\delta_{Y_{\beta}}$ ,  $\beta \in \Gamma_{\alpha}^{w}$ , are independent of  $I_{\alpha}$  and  $I_{\alpha}\delta_{Y_{\alpha}}$ , respectively.

**Remark 2.4.** In the special case of dissociated indicators Theorem 2.3 (i) coincides with Theorem 2.N in Barbour et al. (1992).  $\Box$ 

**Remark 2.5.** Note that the bounds in Theorem 2.3 (i) and (ii) are equal except for the factor  $(1 - \exp{\{-\lambda\}})/\lambda$ .

## 3 The probability of covering of all particles

Assume  $p_1 \ldots, p_k$  are independently and uniformly distributed particles in the rectangle  $\mathbf{A} \subset \mathbb{R}^2$  and let  $C \in \mathcal{K}$ . In this subsection we will derive the probability

$$P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x) \tag{3.1}$$

by means of results in integral geometry, and give some historical background.

First the case of two particles will be considered, so as to give some understanding of the formula for the probability (3.1), given in Theorem 3.2. In this case the probability can be found by fixing one of the particles in an arbitrary position, and considering all possible positions for the other one, such that both particles are covered by some translate of C. By the independence and the uniform distribution of  $p_1$  and  $p_2$ , and the torus convention, it does not matter where the first particle lies, and the quotient of the "possible" area for the second particle and  $\mu(\mathbf{A})$  equals  $P(\exists x \in \mathbf{A} : p_1, p_2 \in C + x)$ .

**Example 3.1.** If  $C \in \mathcal{K}$  is a circle of radius r, the possible area for the second particle is a circle of radius 2r. Then

$$P(\exists x \in \mathbf{A} : p_1, p_2 \in C + x) = 4\pi r^2 / \mu(\mathbf{A}) = 4\mu(C) / \mu(\mathbf{A}).$$

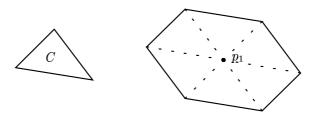


Figure 1: The possible area for the second particle.

**Example 3.2.** Let  $C \in \mathcal{K}$  be a triangle. Figure 1 shows C and the possible area for the second particle, given the position of  $p_1$ . As seen by the figure this area is six times as large as that of the original triangle. Hence

$$P(\exists x \in \mathbf{A} : p_1, p_2 \in C + x) = 6\mu(C)/\mu(\mathbf{A}).$$

By these examples (which will turn out to be extreme, see Corollary 3.3) we learn that for the probability of covering the particles, it is not only the area of the sets which is of importance.

To handle the case of a general  $C \in \mathcal{K}$ , the following simple lemma is of considerable use, by giving an equivalent way of viewing the problem. Recall the definition of the reflection of C at the origin;  $\check{C} = \{-x : x \in C\}$ .

**Lemma 3.1** Let  $x_1, \ldots, x_k$  be arbitrary fixed particles in  $\mathbb{R}^d$  and  $C \subset \mathbb{R}^d$ . Then

$$\exists x \in \mathbb{R}^d \text{ such that } x_1, \dots, x_k \in C + x$$

if and only if

$$\cap_{i=1}^k (\check{C} + x_i) \neq \emptyset.$$

**Proof.** First note that

$$x \in \check{C} + x_i \Leftrightarrow x - x_i \in \check{C} \Leftrightarrow x_i - x \in C \Leftrightarrow x_i \in C + x$$

for every  $i = 1, \ldots, k$ . Hence

$$\bigcap_{i=1}^{k} (\check{C} + x_i) \neq \emptyset \quad \Leftrightarrow \quad \exists x \in \mathbb{R}^d : x \in \bigcap_{i=1}^{k} (\check{C} + x_i) 
\Leftrightarrow \quad \exists x \in \mathbb{R}^d : x_1, \dots, x_k \in C + x.$$

By this lemma it follows that

$$P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x) = P(\cap_{i=1}^k (C + p_i) \neq \emptyset)$$

when the torus convention is used. It is now easy to derive an expression for the probability in case of two particles and a general convex set  $C \in \mathcal{K}$ . Since we may let  $p_1 = O$ , we are looking for the probability that  $p_2$  belongs to the set  $\{x : C \cap (C + x) \neq \emptyset\}$  which is precisely  $C + \check{C}$  by (2.1). Thus

$$P(\exists x \in \mathbf{A} : p_1, p_2 \in C + x) = P((C + p_1) \cap (C + p_2) \neq \emptyset)$$

$$= \mu(C + \check{C})/\mu(\mathbf{A})$$

$$= 2(\mu(C) + \nu(C, \check{C}))/\mu(\mathbf{A}),$$

where the last equality follows from (2.2). What determines the probability is hence the area and the mixed area of the set, where the latter is dependent on the shape of the set. This carries over to the case of more than two particles, as can be seen in Theorem 3.2.

A generalization of the problem formulated in this latter way with non-intersecting sets, is to let the translated sets be of unequal size and shape. In this case there is obviously no dual problem in the context of covering particles. In fact it needs no more effort to handle this generalized case; thus we search the probability

$$P((C_1+p_1)\cap (C_2+p_2)\cap\ldots\cap (C_k+p_k)\neq\emptyset),$$

where  $C_1, \ldots, C_k \in \mathcal{K}$ . Furthermore, for a while we extend the discussion to concern an arbitrary dimension,  $\mathbb{R}^d$ .

A consequence of using the torus convention is that we may assume that the first particle lies at the origin. Because of the assumptions that the origin lies in the centre of **A** and is an interior point of  $C_1, \ldots, C_k$ , and of the restrictions on the sizes of  $C_1, \ldots, C_k$ , all vectors  $(x_2, \ldots, x_k)$  such that

$$C_1 \cap (C_2 + x_2) \cap \ldots \cap (C_k + x_k) \neq \emptyset$$

satisfy  $x_i \in \mathbf{A}$ , i = 2, ..., k, without using the torus convention. Hence, once the assumption that the first particle lies at the origin is made, we may treat  $\mathbf{A}$  as a "normal" d-dimensional rectangle.

Since  $p_2, \ldots, p_k$  are independently and uniformly distributed in  $\mathbf{A}$ , the vector  $(p_2, \ldots, p_k)$  is uniformly distributed in the product space  $\mathbf{A}^{k-1}$ . Let  $\mu_d^{k-1}$  denote the (k-1)-fold product measure of  $\mu_d$ , the d-dimensional Lebesgue measure. Now  $\mu_d^{k-1}(\mathbf{A}^{k-1}) = \mu_d(\mathbf{A})^{k-1}$ , and

$$P(C_1 \cap (C_2 + p_2) \cap \ldots \cap (C_k + p_k) \neq \emptyset) = \mu_d^{k-1} \{(x_2, \ldots, x_k) : x_i \in \mathbb{R}^d, C_1 \cap (C_2 + x_2) \cap \ldots \cap (C_k + x_k) \neq \emptyset\} / \mu_d(\mathbf{A})^{k-1}.$$

To see the connection with integral geometry, we write

$$\mu_d^{k-1}\{(x_2, \dots, x_k) : x_i \in \mathbb{R}^d, C_1 \cap (C_2 + x_2) \cap \dots \cap (C_k + x_k) \neq \emptyset\}$$

$$= \int_{\mathbb{R}^d} \dots \int_{\mathbb{R}^d} V_0(C_1 \cap (C_2 + x_2) \cap \dots \cap (C_k + x_k)) dx_2 \dots dx_k, \tag{3.2}$$

where

$$V_0(C) = \begin{cases} 1 & \text{if } C \neq \emptyset, \\ 0 & \text{if } C = \emptyset. \end{cases}$$

The functional  $V_0(C)$  is one of the so-called *intrinsic volumes* of C,  $V_i(C)$ , i = 0, ..., d, defined for compact, convex subsets of  $\mathbb{R}^d$ . The intrinsic volumes can be defined by the classical Steiner formula as follows. Let  $\mathbf{B}^d$  be the d-dimensional unit ball,  $\kappa_d$  its volume, and  $\lambda \in \mathbb{R}^+$ . Then

$$\mu_d(C + \lambda \mathbf{B}^d) = \sum_{i=0}^d \kappa_{d-i} \lambda^{d-i} V_i(C).$$

The most interesting, and therefore most studied, cases are i = 0, d - 1, d. As already mentioned,  $V_0$  is the indicator of non-empty sets, while  $2V_{d-1}$  is surface area, and  $V_d$  is volume.

As early as 1937 explicit expressions for (3.2) was given for k=2,3 in two and three dimensions. Blaschke (1937) discusses both dimensions while Berwald and Varga (1937) handles three dimensions. In a probabilistic context the planar case including an iterated version (i.e. for an arbitrary number of sets), was rediscovered by Miles (1974), and the two- and three-dimensional cases by Månsson (1996).

In an arbitrary dimension and for general  $V_i$ , the integral in (3.2) is handled in Weil (1990), where it is a special case of an even more general situation. Hence the formula for  $P(\cap_{i=1}^k (C_i + p_i) \neq \emptyset)$  follows directly in an arbitrary dimension. However, in higher dimensions the formulas involve complicated functionals for which explicit descriptions are known only in special cases. Since the setting in this paper is two-dimensional, we present only the probability in this case here.

**Theorem 3.2** Suppose  $C_i \in \mathcal{K}$ , i = 1, ..., k, and that  $p_i$ , i = 1, ..., k, are independently and uniformly distributed particles in  $\mathbf{A}$ , and k = 2, 3, ... Then, using the torus convention,

$$P(\cap_{i=1}^{k}(C_{i}+p_{i})\neq\emptyset) = \left(\sum_{i=1}^{k}\prod_{\substack{j=1\\j\neq i}}^{k}\mu(C_{j}) + \sum_{\substack{i,j=1\\i\neq j}}^{k}\nu(C_{i},\check{C}_{j})\prod_{\substack{l=1\\l\neq i,j}}^{k}\mu(C_{l})\right)\frac{1}{\mu(\mathbf{A})^{k-1}}.$$

In particular, if  $C_i = C$ , i = 1, ..., k, then

$$P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x) = P(\bigcap_{i=1}^k (C + p_i) \neq \emptyset)$$
  
=  $\left(k + k(k-1) \frac{\nu(\check{C}, C)}{\mu(C)}\right) \frac{\mu(C)^{k-1}}{\mu(\mathbf{A})^{k-1}}.$ 

It is the latter part of this theorem which will be useful in the rest of this paper. The following corollary follows directly from Theorem 3.2 and (2.3).

Corollary 3.3 Under the assumptions of Theorem 3.2

$$k^2 \frac{\mu(C)^{k-1}}{\mu(\mathbf{A})^{k-1}} \le P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x) \le (2k^2 - k) \frac{\mu(C)^{k-1}}{\mu(\mathbf{A})^{k-1}},$$

where there is equality on the left if and only if C is centrally symmetric and on the right if and only if C is a triangle.

In this paper the discussions will be carried on in terms of covering particles. But throughout the paper we shall bear in mind that when reading for instance "k uniformly distributed particles are covered by C" we equally well can read "k uniformly translated copies of  $\check{C}$  have a non-empty intersection".

## 4 Poisson approximation

In this section we will consider Poisson approximation of the number of k-subsets of particles which are covered by translates of C, and Poisson process approximation of some point processes which arise in this connection. Results concerning one value of k in the variable and process cases are presented in Subsection 4.1.1 and 4.1.2, respectively, while Subsection 4.1.3 handles joint distributions. The proofs are deferred to Subsection 4.2. Without loss of generality, we will henceforth let  $\mu(\mathbf{A}) = 1$ .

#### 4.1 Main results

#### 4.1.1 The univariate case

As noted in the introduction, the number of k-subsets which are covered by some translate of C can be written as

$$W = \sum_{i=1}^{\binom{n}{k}} I_i, \tag{4.1}$$

where

 $I_i = \left\{ \begin{array}{ll} 1 & \text{if there exists } x \in \mathbf{A} \text{ such that the } i\text{th } k\text{-subset is covered by } C + x, \\ 0 & \text{otherwise.} \end{array} \right.$ 

From Theorem 3.2 we know that

$$E[I_i] = P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x) = \left(k + k(k-1)\frac{\nu(\check{C}, C)}{\mu(C)}\right)\mu(C)^{k-1}, \tag{4.2}$$

 $i=1,\ldots,\binom{n}{k}$ , and hence that

$$\lambda = E[W] = \binom{n}{k} \left( k + k(k-1) \frac{\nu(\check{C}, C)}{\mu(C)} \right) \mu(C)^{k-1}. \tag{4.3}$$

W is a sum of indicators, where those pairs of indicators which concern k-subsets with common particles are dependent while those with no particles in common are independent. In a situation such as this, the local version of the Stein-Chen method is a suitable mean to get a bound on the total variation distance between the distribution of W and a Poisson variable with parameter  $\lambda$ , and it leads to the following theorem.

**Theorem 4.1** Let W and  $\lambda$  be defined by (4.1) and (4.3), respectively. Then

$$d_{TV}(\mathcal{L}(W), \operatorname{Po}(\lambda)) \leq \left\{ \frac{\lambda k^2}{n} + \sum_{l=1}^{k-1} \binom{k}{l} \binom{n-k}{k-l} \mu(\check{C} + C)^{k-l} \right\} (1 - e^{-\lambda}).$$

**Remark 4.2.** In case of k=2 the bound can be somewhat improved by means of Theorem 2.0 in Barbour et al. (1992), which concerns families of strongly dissociated indicator variables. The idea is to reduce the number of indices in  $\Gamma_{\alpha}^{s}$ .

#### 4.1.2 The process case

The purpose of this section is to introduce three point processes determined by positions and sizes of the k-subsets which are covered by some C + x, and consider approximation of these processes by Poisson processes. The point processes are defined as follows.

Let the leftmost particles (the lowest of these in case of ambiguity) in the k-subsets which actually are covered by some translate of C constitute the points of the point process  $\Xi_{\mathbf{A}}$  on  $\mathbf{A}$ . The *size* of a k-subset we define to be the smallest s such that for some  $x \in \mathbf{A}$ , sC + x covers the k-subset. If sizes are attached to the points of  $\Xi_{\mathbf{A}}$ , we get a point process on the space  $\mathbf{A} \times [0,1]$ , which we denote by  $\Xi$ . These sizes are identically, but not independently, distributed with distribution function

$$F(y) = P(\exists x \in \mathbf{A} : p_1, \dots, p_k \in yC + x \mid \exists x \in \mathbf{A} : p_1, \dots, p_k \in C + x)$$

$$= \frac{\left(k + k(k-1)\nu(y\check{C}, yC)/\mu(yC)\right)\mu(yC)^{k-1}}{\left(k + k(k-1)\nu(\check{C}, C)/\mu(C)\right)\mu(C)^{k-1}}$$

$$= y^{2(k-1)}, \tag{4.4}$$

 $0 \le y \le 1$ , by (4.2) and since  $\nu(y\check{C}, yC) = y^2\nu(\check{C}, C)$  and  $\mu(yC) = y^2\mu(C)$ . If we drop the positions and just consider the sizes the result is a point process on [0, 1], which we denote by  $\Xi_{[0,1]}$ .

All these three processes can be written as

$$\sum_{i=1}^{\binom{n}{k}} I_i \delta_{Y_i},$$

where the state space for  $\{Y_i\}$ , denoted by  $\mathcal{Y}$ , is  $\mathbf{A}$ , [0,1] or  $\mathbf{A} \times [0,1]$ . We will first consider  $\Xi$ , for which  $\mathcal{Y} = \mathbf{A} \times [0,1]$ , and we need to derive the measure on  $\mathcal{Y}$  introduced in (2.4);

$$\lambda_i(A) = P(I_i = 1, Y_i \in A).$$

The position of a k-subset which is covered by some translate of C, i.e. its leftmost particle, is uniformly distributed on  $\mathbf{A}$ , since the k particles themselves are uniformly and independently

distributed on A, and the torus convention is used. The size of a k-subset which is covered takes its value in [0,1] and has density function

$$2(k-1)y^{2k-3}$$
,

by (4.4). Furthermore, the size is independent of the position of the k-subset. With  $p = E[I_i]$  and  $\lambda$  defined as in (4.2) and (4.3), respectively, the measure  $\lambda_i$  is thus given by

$$d\lambda_i(x,y) = p \frac{1}{\mu(\mathbf{A})} 2(k-1)y^{2k-3} dx dy = p 2(k-1)y^{2k-3} dx dy, \tag{4.5}$$

 $i = 1, \ldots, \binom{n}{k}$ , and we get

$$d\lambda(x,y) = \sum_{i=1}^{\binom{n}{k}} d\lambda_i(x,y) = \binom{n}{k} p2(k-1)y^{2k-3} dx \, dy = \lambda 2(k-1)y^{2k-3} dx \, dy. \tag{4.6}$$

The bound given below on the total variation distance between  $\mathcal{L}(\Xi)$  and a Poisson process with intensity  $\lambda$  equals the bound on the distance between  $\mathcal{L}(W)$  and a Poisson variable with parameter  $\lambda$  given in Theorem 4.1 if  $\lambda \leq 1$ . If  $\lambda > 1$  it is unfortunately not as good in the process case.

**Theorem 4.3** Let  $\Xi = \sum_{i=1}^{\binom{n}{k}} I_i \delta_{Y_i}$  be the point process on  $\mathbf{A} \times [0,1]$  defined above and let  $\lambda$  be given by (4.6). Then

$$d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda)) \leq \frac{\lambda^2 k^2}{n} + \lambda \sum_{l=1}^{k-1} \binom{k}{l} \binom{n-k}{k-l} \mu(\check{C} + C)^{k-l}.$$

**Remark 4.4.** Since  $\Xi_{\mathbf{A}}$  and  $\Xi_{[0,1]}$  and the corresponding Poisson processes are obtained as measurable mappings from  $\Xi$  and  $\mathrm{Po}(\lambda)$ , respectively, it follows that

$$d_{TV}(\mathcal{L}(\Xi_{\mathcal{V}}), \text{Po}(\lambda_{\mathcal{V}})) \leq d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda)),$$

where  $\mathcal{Y} = \mathbf{A}$  or [0,1], and  $\lambda_{\mathcal{Y}}$  is the measure corresponding to (4.6). Hence the bound in Theorem 4.3 holds also when these processes are concerned.

#### 4.1.3 The multivariate case

For  $C^{(k_j)} \in \mathcal{K}$ , where  $j = 1, 2, ..., m, k_j \in \{2, 3, ...\}$  and  $k_i \neq k_j$  if  $i \neq j$ , let

$$W^{(k_j)} = \sum_{i=1}^{\binom{n}{k_j}} I_{k_j,i}, \tag{4.7}$$

and

$$\Xi^{(k_j)} = \sum_{i=1}^{\binom{n}{k_j}} I_{k_j,i} \delta_{Y_{k_j,i}}, \tag{4.8}$$

where  $I_{k_j,i}=1$  if the i'th  $k_j$ -subset is covered by some translate of  $C^{(k_j)}$  and 0 otherwise. We will now consider approximation of the distribution of the vectors  $(W^{(k_1)},\ldots,W^{(k_m)})$  and  $(\Xi^{(k_1)},\ldots,\Xi^{(k_m)})$  by the vectors of corresponding independent Poisson variables and Poisson processes, respectively.

From (4.3) we know that

$$\lambda^{(k_j)} = E[W^{(k_j)}] = \binom{n}{k_j} \left( k_j + k_j (k_j - 1) \frac{\nu(\check{C}^{(k_j)}, C^{(k_j)})}{\mu(C^{(k_j)})} \right) \mu(C^{(k_j)})^{k_j - 1}, \tag{4.9}$$

and furthermore, by (4.6),

$$d\lambda^{(k_j)}(x,y) = \sum_{i=1}^{\binom{n}{k_j}} d\lambda_i^{(k_j)}(x,y) = \lambda^{(k_j)} 2(k_j - 1) y^{2k_j - 3} dx dy, \tag{4.10}$$

is the measure on  $\mathbf{A} \times [0,1]$ , which is connected to  $\Xi^{(k_j)}$ .

**Theorem 4.5** Let  $W^{(k_j)}$ ,  $\Xi^{(k_j)}$ ,  $\lambda^{(k_j)}$  and  $\lambda^{(k_j)}$  be defined by (4.7), (4.8), (4.9) and (4.10), respectively. Then

$$\begin{split} d_{TV}(\mathcal{L}(\{W^{(k_{j})}\}_{j=1}^{m}), \Pi_{j=1}^{m} \mathrm{Po}(\lambda^{(k_{j})})) & \leq & d_{TV}(\mathcal{L}(\{\Xi^{(k_{j})}\}_{j=1}^{m}), \Pi_{j=1}^{m} \mathrm{Po}(\pmb{\lambda}^{(k_{j})})) \\ & \leq & \sum_{i=1}^{m} \binom{n}{k_{j}} (p_{k_{j}}^{2} + p_{k_{j}} E[Z_{k_{j},1}] + E[I_{k_{j},1} Z_{k_{j},1}]), \end{split}$$

where

$$E[Z_{k_{j},1}] = \sum_{l=1}^{k_{j}-1} {k_{j} \choose l} {n-k_{j} \choose k_{j}-l} p_{k_{j}} + \sum_{\substack{s=1\\s\neq j}}^{m} \sum_{l=1}^{\min\{k_{j},k_{s}\}} {k_{j} \choose l} {n-k_{j} \choose k_{s}-l} p_{k_{s}}, \qquad (4.11)$$

$$E[Z_{k_{j},1}I_{k_{j},1}] \leq$$

$$\sum_{l=1}^{k_{j}-1} {k_{j} \choose l} {n-k_{j} \choose k_{j}-l} p_{k_{j}} \mu(C^{(k_{j})} + \check{C}^{(k_{j})})^{k_{j}-l}$$

$$+ \sum_{\substack{s=1 \ s \neq i}}^{m} \sum_{l=1}^{\min\{k_{j},k_{s}\}} {k_{j} \choose l} {n-k_{j} \choose k_{s}-l} \min\{p_{k_{j}} \mu(C^{(k_{s})} + \check{C}^{(k_{s})})^{k_{s}-l}, p_{k_{s}} \mu(C^{(k_{j})} + \check{C}^{(k_{j})})^{k_{j}-l}\}$$

$$(4.12)$$

and

$$p_{k_j} = \left(k_j + k_j(k_j - 1) \frac{\nu(C^{(k_j)}, \check{C}^{(k_j)})}{\mu(C^{(k_j)})}\right) \mu(C^{(k_j)})^{k_j - 1}. \tag{4.13}$$

#### 4.2 Proofs

We will start to approximate  $\sum_{j=1}^{m} W^{(k_j)}$  by a Poisson variable with expectation

$$E[\sum_{j=1}^{m} W^{(k_j)}] = \sum_{j=1}^{m} \lambda^{(k_j)},$$

where  $W^{(k_j)}$  and  $\lambda^{(k_j)}$  are defined as in (4.7) and (4.9), respectively. This approximation is no aim in itself but is more a means to be used in the rest of the proofs.

First we must introduce some new, local, notation. Let

$$\begin{array}{lcl} \Gamma^{j}_{j,i} & = & \{l:l=1,\ldots,\binom{n}{k_{j}}, l\neq i, (i\text{th }k_{j}\text{-subset}\cap l\text{th }k_{j}\text{-subset})\neq\emptyset\}, \ j=1,\ldots,m, \\ \Gamma^{j}_{s,i} & = & \{l:l=1,\ldots,\binom{n}{k_{s}}, (i\text{th }k_{j}\text{-subset}\cap l\text{th }k_{s}\text{-subset})\neq\emptyset\}, \\ & & j=1,\ldots,m, \ s=1,\ldots,m, \ s\neq j, \\ Z_{k_{j},i} & = & \sum_{s=1}^{m}\sum_{l\in\Gamma^{j}_{s,i}}I_{k_{s},l}, \ j=1,\ldots,m, \end{array}$$

so that  $Z_{k_j,i}$  equals the sum of the number of  $k_s$ -subsets which are covered by some  $C^{(k_s)} + x$ ,  $s = 1, \ldots, m$ , and have some particles in common with the *i*th  $k_j$ -subset.

Theorem 4.6 With the notation introduced above

$$d_{TV}(\mathcal{L}(\sum_{j=1}^{m} W^{(k_{j})}), \operatorname{Po}(\sum_{j=1}^{m} \lambda^{(k_{j})})) \leq \left(\sum_{j=1}^{m} \binom{n}{k_{j}} (p_{k_{j}}^{2} + p_{k_{j}} E[Z_{k_{j},1}] + E[I_{k_{j},1} Z_{k_{j},1}])\right) \frac{1 - \exp\{-\sum_{j=1}^{m} \lambda^{(k_{j})}\}}{\sum_{j=1}^{m} \lambda^{(k_{j})}},$$
(4.14)

where  $E[Z_{k_j,1}]$ ,  $E[I_{k_j,1}Z_{k_j,1}]$  and  $p_{k_j}$  are given by (4.11), (4.12) and (4.13), respectively.

**Proof.** Note that if we let  $\Gamma$  consist of the indices of all  $k_j$ -subsets of  $\{1, 2, ..., n\}$ , j = 1, ..., m, and  $\Gamma_{\alpha}^s = \{\beta \in \Gamma : \beta \neq \alpha, \beta \cap \alpha \neq \emptyset\}$ , then we can write

$$\sum_{j=1}^{m} W^{(k_j)} = \sum_{j=1}^{m} \sum_{i=1}^{\binom{n}{k_j}} I_{k_j,i} = \sum_{\alpha \in \Gamma} I_{\alpha},$$

where  $I_{\alpha} = 1$  if all  $p_i$ ,  $i \in \alpha$ , are covered by some  $C^{(|\alpha|)} + x$ ,  $x \in \mathbf{A}$ . Furthermore, if  $\alpha$  is the index of the *i*th  $k_j$ -subset, then

$$Z_{k_j,i} = \sum_{s=1}^m \sum_{i \in \Gamma_{s-1}^j} I_{k_s,i} = \sum_{\beta \in \Gamma_{\alpha}^s} I_{\beta}.$$

Hence it is clear that the assumptions of Theorem 2.3 (i) are satisfied, and (4.14) follows directly. We need now to derive the expression and bound for  $E[Z_{k_j,1}]$  and  $E[I_{k_j,1}Z_{k_j,1}]$ , respectively.

There are  $\binom{k_j}{l}\binom{n-k_j}{k_j-l}$  different  $k_j$ -subsets with exactly l particles in common with the 1st  $k_j$ -subset,  $l=1,\ldots,k_j-1$ , and  $\binom{k_j}{l}\binom{n-k_j}{k_s-l}$  different  $k_s$ -subsets with exactly l particles in common with the 1st  $k_j$ -subset,  $l=1,\ldots,\min\{k_j,k_s\}$ . By changing the index of summation in the right-hand side of

$$E[Z_{k_j,1}] = \sum_{s=1}^{m} \sum_{i \in \Gamma_{s,1}^j} E[I_{k_s,i}] = \sum_{s=1}^{m} \sum_{i \in \Gamma_{s,1}^j} p_{k_j}$$

we immediately get (4.11).

Now, for each  $l=1,\ldots,k_j-1$ , let  $I_{k_j}^l$  be an indicator with the same distribution as a  $k_j$ -subset which has l particles in common with the 1st  $k_j$ -subset, and let  $I_{k_s}^l$  be the corresponding variable but concerning a  $k_s$ -subset which has l particles in common with the 1st  $k_j$ -subset, where  $l=1,\ldots,\min\{k_j,k_s\}$ . Then  $E[Z_{k_j,1}I_{k_j,1}]$  can be rewritten as

$$E[Z_{k_{j},1}I_{k_{j},1}] = \sum_{l=1}^{k_{j}-1} {k_{j} \choose l} {n-k_{j} \choose k_{j}-l} E[I_{k_{j},1}I_{k_{j}}^{l}] + \sum_{\substack{s=1 \ s \neq j}}^{m} \sum_{l=1}^{\min\{k_{j},k_{s}\}} {k_{j} \choose l} {n-k_{j} \choose k_{s}-l} E[I_{k_{j},1}I_{k_{s}}^{l}].$$

$$(4.15)$$

The conditional probability  $P(I_{k_j}^l = 1 \mid I_{k_j,1} = 1)$  is bounded above by the probability that the  $k_j - l$  not common particles lie in the set  $C^{(k_j)} + \check{C}^{(k_j)}$ , centred on any of the l common particles:

$$P(I_{k_j}^l = 1 \mid I_{k_j,1} = 1) \le \mu(C^{(k_j)} + \check{C}^{(k_j)})^{k_j - l},$$

and hence

$$E[I_{k_{j},1}I_{k_{j}}^{l}] = P(I_{k_{j}}^{l} = 1 \mid I_{k_{j},1} = 1)p_{k_{j}}$$

$$\leq \mu(C^{(k_{j})} + \check{C}^{(k_{j})})^{k_{j}-l}p_{k_{j}}. \tag{4.16}$$

As above,

$$E[I_{k_j,1}I_{k_s}^l] \le \mu(C^{(k_s)} + \check{C}^{(k_s)})^{k_s-l}p_{k_j}, \tag{4.17}$$

and by symmetry, another bound for  $E[I_{k_i,1}I_{k_s}^l]$  is

$$E[I_{k_j,1}I_{k_s}^l] \leq \mu(C^{(k_j)} + \check{C}^{(k_j)})^{k_j-l}p_{k_s}. \tag{4.18}$$

Insert (4.16)–(4.18) in (4.15) to obtain (4.12).

**Proof of Theorem 4.1** Letting m=1 and  $k_1=k$  in Theorem 4.6 and noting that  $p_k=\lambda/\binom{n}{k}$  gives

$$d_{TV}(\mathcal{L}(W), \operatorname{Po}(\lambda)) = \left\{ \frac{\lambda}{\binom{n}{k}} + \frac{\lambda}{\binom{n}{k}} \sum_{l=1}^{k-1} \binom{k}{l} \binom{n-k}{k-l} + \sum_{l=1}^{k-1} \binom{k}{l} \binom{n-k}{k-l} \mu(\check{C} + C)^{k-l} \right\} (1 - e^{-\lambda}).$$

$$(4.19)$$

Let X be a hypergeometric random variable such that

$$P(X = l) = \frac{\binom{k}{l} \binom{n-k}{k-l}}{\binom{n}{k}}, \quad l = 0, 1, \dots, k, \quad \text{and} \quad E[X] = \frac{k^2}{n}.$$

Then

$$\frac{1}{\binom{n}{k}} \sum_{l=1}^{k-1} \binom{k}{l} \binom{n-k}{k-l} = P(X \ge 1) - \frac{1}{\binom{n}{k}} \le \frac{k^2}{n} - \frac{1}{\binom{n}{k}}$$

by Markov's inequality, and the first two terms in the major brackets of (4.19) can be bounded by  $\lambda k^2/n$ , by which the theorem is proved.

**Remark 4.7.** Theorem 4.1 could be derived more directly by using (3.2) p. 35 in Barbour et al. (1992).

**Proof of Theorem 4.3.** Positions pertaining to k-subsets with common particles are dependent, otherwise they are independent. The same holds for the sizes. Furthermore this dependence structure is the same as that of  $\{I_i: i=1,\ldots,\binom{n}{k}\}$ . Hence the family  $\{I_i\delta_{Y_i}: i=1,\ldots,\binom{n}{k}\}$  satisfies the assumptions of Theorem 2.3 (ii). As noted in Remark 2.5, in such a case the bound of  $d_{TV}(\mathcal{L}(\Xi), \operatorname{Po}(\lambda))$  given in Theorem 2.3 (ii) coincides with the bound of  $d_{TV}(\mathcal{L}(W), \operatorname{Po}(\lambda))$  given in Theorem 2.3 (i), except for the factor  $(1-\exp(\lambda))/\lambda$ . We used Theorem 4.6 in the special case where j=1 to bound  $d_{TV}(\mathcal{L}(W), \operatorname{Po}(\lambda))$ . However, in that case Theorem 4.6 and Theorem 2.3 (i) coincide, and the result follows without further calculations.

**Proof of Theorem 4.5.** To prove this theorem, we use another point process  $\widetilde{\Xi}$  which is the sum of the processes  $\Xi^{(k_j)}$ ,  $j=1,\ldots,m,\,k_j\in\{2,3,\ldots\}$ , but with the additional information of which

process  $\Xi^{(k_j)}$  each point originates from. Hence, if  $\Gamma^{(k_j)}$  contains the indices of the  $k_j$ -subsets, and  $\Gamma = \bigcup_{j=1}^m \Gamma^{(k_j)}$ , let

$$\widetilde{\Xi} = \sum_{\alpha \in \Gamma} I_{\alpha} \delta_{\widetilde{Y}_{\alpha}}$$

where the  $\widetilde{Y}_{\alpha}$ 's take their values in

$$\widetilde{\mathcal{Y}} = \mathbf{A} \times [0,1] \times \{k_1, k_2, \dots, k_m\}.$$

The first step is to derive  $\widetilde{\lambda}_{\alpha}(A) = P(I_{\alpha} = 1, \widetilde{Y}_{\alpha} \in A)$ , where  $A \subset \widetilde{\mathcal{Y}}$ . Note that if  $\alpha \in \Gamma^{(k_j)}$ , then

$$d\widetilde{\lambda}_{\alpha}(x,y,k_j) = p_{k_j} 2(k_j - 1) y^{2k_j - 3} dx dy,$$

by (4.5), and  $d\widetilde{\lambda}_{\alpha}(x,y,z) = 0$  for  $z \neq k_j$ . Hence, for  $(x,y,z) \in \widetilde{\mathcal{Y}}$ , let

$$d\widetilde{\lambda}_{\alpha}(x,y,z) = p_z 2(z-1)y^{2z-3} 1\{\alpha \in \Gamma^{(z)}\} dx dy,$$

so that

$$d\widetilde{\lambda}(x,y,z) = \sum_{\alpha \in \Gamma} d\widetilde{\lambda}_{\alpha}(x,y,z) = \sum_{\alpha \in \Gamma^{(z)}} d\widetilde{\lambda}_{\alpha}(x,y,z) = \lambda^{(z)} 2(z-1) y^{2z-3} dx dy.$$

Since

$$W^{(k_j)} = \int_{\mathbf{A}} \int_{[0,1]} \Xi^{(k_j)} \{ dx, dy \}$$

is a functional of  $\Xi^{(k_j)}$ , and

$$\Xi^{(k_j)}\{dx, dy\} = \widetilde{\Xi}\{dx, dy, k_j\}$$

in turn is a measurable mapping of  $\widetilde{\Xi}$ , it follows that

$$d_{TV}(\mathcal{L}(\{W^{(k_j)}\}_{j=1}^m), \Pi_{j=1}^m \text{Po}(\lambda^{(k_j)})) \leq d_{TV}(\mathcal{L}(\{\Xi^{(k_j)}\}_{j=1}^m), \Pi_{j=1}^m \text{Po}(\lambda^{(k_j)})) \\ < d_{TV}(\mathcal{L}(\widetilde{\Xi}), \text{Po}(\widetilde{\lambda})).$$

For a fixed k we could use the bound on  $d_{TV}(\mathcal{L}(W), \text{Po}(\lambda))$  to get a bound on  $d_{TV}(\mathcal{L}(\Xi), \text{Po}(\lambda))$ , because of the dependence structure of  $\{I_i\delta_{Y_i}, i=1,\ldots,\binom{n}{k}\}$ . Similarly we get a bound on  $d_{TV}(\mathcal{L}(\widetilde{\Xi}), \text{Po}(\widetilde{\lambda}))$  by means of the bound on  $d_{TV}(\mathcal{L}(\sum_{j=1}^m W^{(k_j)}), \text{Po}(\sum_{j=1}^m \lambda^{(k_j)}))$  given in Theorem 4.6.

# 5 Asymptotic results

For a fixed C and k, the errors made by Poisson approximating of course get large when n does. We will now consider asymptotic aspects of the previous results as k is kept fixed, but the areas of the sets decrease as n grows. For instance how fast the areas of sets in a sequence must decrease for the total variation distances to go to zero, if these rates are different in the uni- and multivariate cases, and if there is any difference between the variable and the process cases. Since n is no longer fixed, the quantities which change with n are here indexed by n.

#### 5.1 The univariate case

Let  $\{C_n\}_{n=1}^{\infty}$ ,  $C_n \in \mathcal{K}$ , be a sequence of sets. To get a bound on the variation distance  $d_{TV}(\mathcal{L}(W_n), \text{Po}(\lambda_n))$ , which is given in Theorem 4.1, valid for all shapes of the set  $C_n$  we use that

$$\mu(\check{C}_n + C_n) \le 6\mu(C_n)$$

by (2.2) and (2.3). Furthermore, the expectation of  $W_n$ , given in (4.3), can be bounded by

$$\lambda_n \le \frac{n^k (2k^2 - k)\mu(C_n)^{k-1}}{k!},$$

by Corollary 3.3. These bounds, together with  $(1 - e^{-\lambda_n}) \leq \min(1, \lambda_n)$ , inserted in the bound in Theorem 4.1 yields

$$d_{TV}(\mathcal{L}(W_n), \text{Po}(\lambda_n)) \leq \left\{ \frac{(2k^2 - k)k(n\mu(C_n))^{k-1}}{(k-1)!} + \sum_{l=1}^{k-1} \binom{k}{l} \frac{(6n\mu(C_n))^{k-l}}{(k-l)!} \right\} \times \min \left\{ 1, \frac{n^k(2k^2 - k)\mu(C_n)^{k-1}}{k!} \right\},$$
(5.1)

valid for all  $C_n \in \mathcal{K}$ .

**Theorem 5.1** Let  $W_n$  and  $\lambda_n$  be defined as in (4.1) and (4.3), respectively.

(i) For any sequence of sets,  $\{C_n\}_{n=1}^{\infty}$ ,  $C_n \in \mathcal{K}$ ,

$$d_{TV}(\mathcal{L}(W_n), \text{Po}(\lambda_n)) = O\left(\min\{1, n^k \mu(C_n)^{k-1}\} \sum_{l=1}^{k-1} (n\mu(C_n))^{k-l}\right).$$
 (5.2)

(ii) For a sequence of sets  $\{C_n\}_{n=1}^{\infty}$ ,  $C_n \in \mathcal{K}$ , with  $\mu(C_n) = o(n^{-t})$ , where  $t \geq 1$  is constant, the bound tends to zero and is of the order

$$d_{TV}(\mathcal{L}(W_n), \text{Po}(\lambda_n)) = \begin{cases} o(n^{1-t}) & \text{if } 1 \le t \le k/(k-1), \\ o(n^{k(1-t)+1}) & \text{if } t > k/(k-1). \end{cases}$$
(5.3)

The conditions t > k/(k-1) and  $t \le k/(k-1)$  are here equivalent to  $\lambda_n \to 0$  and  $\lambda_n \to \infty$ , respectively.

**Proof.** (i) Follows directly from (5.1).

(ii) Since  $t \geq 1$ , it follows that  $n\mu(C_n) < 1$  asymptotically. Hence it is the (k-1)st term (i.e. l = k-1) in the sum in (5.2) which is dominating, and it is of order  $o(n^{1-t})$ . Furthermore

$$\min\{1, n^k \mu(C_n)^{k-1}\} = \left\{ \begin{array}{ll} o(1) & \text{if } 1 \leq t \leq k/(k-1), \\ o(n^{k(1-t)+t}) & \text{if } t > k/(k-1). \end{array} \right.$$

Combining these facts yields (5.3). The last statement follows from the fact that

$$\lambda_n = O(n^k \mu(C_n)^{k-1}) = o(n^{k-t(k-1)}).$$

**Remark 5.2.** Note that it is the (k-1)st term of the sum which determines the rates in the second part of the theorem. This term concerns the dependence between two indicators connected to k-subsets with k-1 common particles.

**Remark 5.3.** The condition  $\mu(C_n) = o(n^{-1})$  is equivalent to that  $\lambda_n = o(n)$ . Hence the expected number of k-subsets which are covered may tend to infinity at the rate o(n) and still the variation distance tends to zero.

Corollary 5.4 Let  $\{C_n\}_{n=1}^{\infty}$  be a sequence of sets such that  $C_n \in \mathcal{K}$  and

$$E[W_n] \to \lambda \quad as \quad n \to \infty,$$

where  $0 \le \lambda < \infty$ , and let  $\mathcal{L}(X) = Po(\lambda)$ . Then

$$W_n \stackrel{\mathcal{D}}{\to} X$$
 as  $n \to \infty$ .

**Proof.** Since

$$E[W_n] = \binom{n}{k} \left( k + k(k-1)\nu(C_n, \check{C}_n) / \mu(C_n) \right) \mu(C_n)^{k-1} \to \lambda \quad \text{as } n \to \infty,$$

and  $\nu(C_n, \check{C}_n)/\mu(C_n) \geq 1$ , for all n, by (2.3), there exists  $0 \leq c < \infty$  such that

$$\mu(C_n) < cn^{-\frac{k}{k-1}}.$$

By Theorem 5.1 (ii) the total variation distance between  $\mathcal{L}(W_n)$  and  $Po(\lambda_n)$  tends to zero in this case, which combined with the fact that

$$d_{TV}(\operatorname{Po}(\lambda_n), \operatorname{Po}(\lambda)) \to 0$$

if  $\lambda_n \to \lambda$ , implies convergence in distribution by the triangle inequality.

This subsection is concluded by an application of Theorem 5.1. Let  $M_{C_n}$  be the maximal number of n independent and uniformly distributed particles on  $\mathbf{A}$  which are covered by  $C_n + x$  for some  $x \in \mathbf{A}$ . We will study the distribution of  $M_{C_n}$  for different sequences of sets, and how its asymptotic distribution depends on the chosen sequence.

Suppose c>0 and  $t\in\mathbb{R}$  are constants and let  $\{C_n\}_{n=1}^{\infty}$  be a sequence of subsets of  $\mathcal{K}$  such that

$$\mu(C_n) = cn^{-t}.$$

Then the expectation of the number of k-subsets which are covered by some translate of  $C_n$  is

$$\lambda_n = E[W_n] = \binom{n}{k} \left( k + k(k-1) \frac{\nu(C_n, \check{C}_n)}{\mu(C_n)} \right) c^{k-1} n^{-t(k-1)}.$$
 (5.4)

The limit of this expectation depends on the value of t:

$$\lambda_n \to \begin{cases} \infty & \text{if } t < k/(k-1) \\ 0 & \text{if } t > k/(k-1) \end{cases} \text{ as } n \to \infty,$$
 (5.5)

and if t = k/(k-1)

$$\frac{k^2c^{k-1}}{k!} \le \liminf_{n \to \infty} \lambda_n \le \limsup_{n \to \infty} \lambda_n \le \frac{(2k^2 - k)c^{k-1}}{k!},$$

where the bounds follow by Corollary 3.3.

**Theorem 5.5** Suppose that  $k \in \{2,3,\ldots\}$ , c > 0 and let  $\{C_n\}_1^{\infty}$  be a sequence of subsets of K with

$$\mu(C_n) = cn^{-t}.$$

If t = k/(k-1) and  $\lambda_n \to \lambda$  as  $n \to \infty$ , then

$$P(M_{C_n} = i) \to \begin{cases} e^{-\lambda} & i = k - 1\\ 1 - e^{-\lambda} & i = k\\ 0 & i \neq k - 1, k \end{cases} \quad as \ n \to \infty.$$

If 
$$(k+1)/k < t < k/(k-1)$$
, then

$$P(M_{C_n} = i) \to \begin{cases} 1 & i = k \\ 0 & i \neq k \end{cases} \quad as \ n \to \infty.$$

**Remark 5.6.** If all  $C_n$  are of the same shape and t = k/(k-1), then  $\lambda_n \to \lambda$ , since for some  $1 \le a \le 2$ ,  $\nu(\check{C}_n, C_n)/\mu(C_n) = a$  for all n.

**Proof.** First we let t = k/(k-1). It then follows from Corollary 5.4 that

$$P(M_{C_n} < k) = P(W_n = 0) \quad \to \quad \exp\{-\lambda\} \tag{5.6}$$

as  $n \to \infty$ .

Let  $W_{n,k-1}$  and  $W_{n,k+1}$  denote the number of covered (k-1)- and (k+1)-subsets, respectively, and let  $\lambda_{n,k-1}$  and  $\lambda_{n,k+1}$  be their expectations. If k=2 then  $\lambda_{n,k-1} \to \infty$  since every single particle is of course covered by some  $C_n + x$ . Furthermore, since t < (k-1)/(k-2), k > 2, and t > (k+1)/k we get by (5.5) that

$$\lambda_{n,k-1} \to \infty$$
 and  $\lambda_{n,k+1} \to 0$  as  $n \to \infty$ . (5.7)

Hence the probability that  $M_{C_n}$  is strictly less than k-1 approaches 0 since  $\mu(C_n) = cn^{-k/(k-1)}$ , so by Theorem 5.1 (ii)

$$P(M_{C_n} < k - 1) = P(W_{n,k-1} = 0)$$
  
=  $\exp\{-\lambda_{n,k-1}\} + o(n^{-\frac{1}{k-1}}) \to 0$  as  $n \to \infty$ ,

which together with (5.6) yields

$$P(M_{C_n} = k - 1) \to \exp\{-\lambda\}.$$

Next we consider the probability that  $M_{C_n}$  is greater than k:

$$P(M_{C_n} \ge k+1) = P(W_{n,k+1} \ge 1)$$
 
$$\le E[W_{n,k+1}] \to 0 \quad \text{as } n \to \infty ,$$

by Markov's inequality and (5.7). The only possibility that remains for  $M_{C_n}$  in the limit is to be equal to k:

$$P(M_{C_n} = k) = 1 - P(M_{C_n} < k) - P(M_{C_n} > k) \to 1 - \exp\{-\lambda\}$$

as  $n \to \infty$ . Hence the theorem is proved for t = k/(k-1).

For (k+1)/k < t < k/(k-1) it can be shown that  $P(M_{C_n} < k) \to 0$  and  $P(M_{C_n} \ge k+1) \to 0$  by arguments similar to those above. Hence

$$P(M_{C_n} = k) = 1 - P(M_{C_n} < k) - P(M_{C_n} > k + 1) \to 1$$
 as  $n \to \infty$ 

and the theorem is proved.

#### 5.2 The process case

Let  $\{\Xi_n\}_{n=1}^{\infty}$  and  $\{\lambda_n\}_{n=1}^{\infty}$  be the sequences of processes and measures, respectively, introduced in Subection 4.1.2, which correspond to the sequence of sets  $\{C_n\}_{n=1}^{\infty}$ . The bound on  $d_{TV}(\mathcal{L}(\Xi_n), \operatorname{Po}(\lambda_n))$  in Theorem 4.3 equals the bound on  $d_{TV}(\mathcal{L}(W_n), \operatorname{Po}(\lambda_n))$  in Theorem 4.1, except that it lacks the factor  $\min(1, 1/\lambda_n)$ , and is hence not as good when  $\lambda_n > 1$ . To obtain convergence of  $d_{TV}(\mathcal{L}(\Xi_n), \operatorname{Po}(\lambda_n))$  to zero, we can therefore not allow the areas of the sets to decrease as slowly as in the previous case of sequences of variables. The order terms in Theorem 5.1 valid for  $\lambda_n \leq 1$ , i.e. when  $\min\{1, n^k\mu(C_n)^{k-1}\} = n^k\mu(C_n)^{k-1}$  in (i), and when  $t \geq k/(k-1)$  in (ii), will be valid for all  $\lambda_n$  here, and the counterpart of Theorem 5.1 then reads:

**Theorem 5.7** Let  $\Xi_n$  and  $\lambda_n$  be defined as in Theorem 4.3. Then, for any sequence of sets  $\{C_n\}_{n=1}^{\infty}, C_n \in \mathcal{K}$ ,

$$d_{TV}(\mathcal{L}(\Xi_n), \text{Po}(\lambda_n)) = O(\sum_{l=1}^{k-1} n^{2k-l} \mu(C_n)^{2k-l-1}).$$
(5.8)

For a sequence of sets  $\{C_n\}_{n=1}^{\infty}$ ,  $C_n \in \mathcal{K}$ , with  $\mu(C_n) = o(n^{-t})$ , where  $t \geq (k+1)/k$  is constant, this bound tends to zero and is of the order

$$d_{TV}(\mathcal{L}(\Xi_n), \text{Po}(\lambda_n)) = o(n^{k(1-t)+1}).$$

Note that the condition for convergence in the variable case,  $\mu(C_n) = o(1/n)$ , does not depend on k. Here the condition is  $\mu(C_n) = o(n^{-(k+1)/k})$ ; the smaller k the faster must the areas decrease.

#### 5.3 The multivariate case

By studying the bound on the total variation distance in the multivariate case, given in Theorem 4.5, we find that it is  $\binom{n}{k_j}E[I_{k_j,1}Z_{k_j,1}]$  which has to be examined to achieve conditions for convergence. Note that it is obvious that we need at least that

$$n\mu(C_n^{(k_j)}) \to 0, \tag{5.9}$$

which was the condition for convergence in the univariate case (Theorem 5.1 (ii)). Hence it is enough to consider the terms corresponding to  $l=k_j-1$  and  $l=\min\{k_j,k_s\}$  of the bound on  $E[I_{k_j,1}Z_{k_j,1}]$  given in (4.12), since they are dominating if (5.9) is fulfilled. Collect the n-dependent parts of these terms, to get the following conditions on  $\mu(C_n^{(k_j)})$  for convergence of  $d_{TV}(\mathcal{L}(\{\Xi_n^{(k_j)}\}_{j=1}^m), \Pi_{j=1}^m \text{Po}(\lambda_n^{(k_j)}))$  to zero:

$$n^{k_j+1}\mu(C_n^{(k_j)})^{k_j} \to 0 \tag{5.10}$$

and

$$m^{k_{s}+k_{j}-\min\{k_{j},k_{s}\}} \times \\ \min \left\{ \mu(C_{n}^{(k_{s})})^{k_{s}-1} \mu(C_{n}^{(k_{j})})^{k_{j}-\min\{k_{j},k_{s}\}}, \mu(C_{n}^{(k_{j})})^{k_{j}-1} \mu(C_{n}^{(k_{s})})^{k_{s}-\min\{k_{j},k_{s}\}} \right\} \\ \to 0$$

$$(5.11)$$

as  $n \to \infty$ , for all  $j, s = 1, ..., m, j \neq s$ .

In the previous case, concerning one point process, it was the  $(k_j-1)$ st term in the sum in (5.8) which decided at which rate the areas of the sets in the sequence  $\{C_n^{(k_j)}\}$  must decrease for the distance to tend to zero. The important part of this term originates from the expectation of the product of two indicators connected to  $k_j$ -subsets with  $k_j-1$  common particles (i.e.  $E[I_{k_j,1}I_{k_j}^{k_j-1}]$  in (4.15), using the notation in the proof of Theorem 4.6). This expectation is of importance also in this multivariate case and turns up as condition (5.10). The other condition, (5.11), comes from the expectation of products of indicators pertaining to  $k_j$ -subsets and  $k_s$ -subsets with min $\{k_j, k_s\}$  particles in common  $(E[I_{k_s,1}I_{k_j}^{\min\{k_j,k_s\}}])$ .

If  $k_j < k_s$  condition (5.11) reads

$$\min \left\{ n^{k_s} \mu(C_n^{(k_s)})^{k_s - 1}, n^{k_s} \mu(C_n^{(k_j)})^{k_j - 1} \mu(C_n^{(k_s)})^{k_s - k_j} \right\} \to 0 \quad \text{as} \quad n \to \infty,$$

and we can state the following theorem.

**Theorem 5.8** If  $\{C_n^{(k_j)}\}_{n=1}^{\infty}$ ,  $j=1,\ldots,m,\ k_j\in\{2,3,\ldots\}$ , are sequences of sets such that

$$n^{k_j+1}\mu(C_n^{(k_j)})^{k_j} \to 0 \quad \text{as } n \to \infty,$$
 (5.12)

for all j = 1, ..., m, and

$$\min \left\{ n^{k_s} \mu(C_n^{(k_s)})^{k_s - 1}, n^{k_s} \mu(C_n^{(k_j)})^{k_j - 1} \mu(C_n^{(k_s)})^{k_s - k_j} \right\} \to 0 \quad \text{as } n \to \infty,$$
 (5.13)

for all j, s = 1, ..., m, such that  $k_j < k_s$ , then

$$d_{TV}\left(\mathcal{L}(\{W_n^{(k_j)}\}_{j=1}^m), \Pi_{j=1}^m \text{Po}(\lambda_n^{(k_j)})\right) \le d_{TV}\left(\mathcal{L}(\{\Xi_n^{(k_j)}\}_{j=1}^m), \Pi_{j=1}^m \text{Po}(\lambda_n^{(k_j)})\right) \to 0$$

as  $n \to \infty$ .

As expected, sequences  $\{C_n^{(k_j)}\}$  with  $\mu(C_n^{(k_j)}) = o(n^{-(k_j+1)/k_j})$ , j = 1, ..., m, which was the critical rate for convergence in total variation norm in case of one process, are not decreasing fast enough here; (5.12) is fulfilled, but (5.13) is not. An example which satisfies both conditions is

$$\mu(C_n^{(k_j)}) = o(n^{-\frac{k_j+1}{k_j} - \frac{1}{k_j^3 - 1}}),$$

 $j = 1, \ldots, m$ , while

$$\mu(C_n^{(k_j)}) = o(n^{-\frac{k_j+1}{k_j} - \frac{1}{k_j^3}}),$$

 $j = 1, \dots, m$ , do not decrease fast enough.

If the expected numbers of covered  $k_j$ -subsets,  $W_n^{(k_j)}$ ,  $j=1,\ldots,m$ , converge to finite values, then joint convergence in distribution follows immediately.

Corollary 5.9 If  $E[W_n^{(k_j)}] \to \lambda^{(k_j)}$ ,  $0 \le \lambda^{(k_j)} < \infty$ , j = 1, ..., m, and  $\mathcal{L}(Z^{(k_j)}) = \text{Po}(\lambda^{(k_j)})$  are independent, then

$$(W_n^{(k_1)},\dots,W_n^{(k_m)}) \overset{\mathcal{D}}{\to} (Z^{(k_1)},\dots,Z^{(k_m)}) \qquad \text{as} \quad n \to \infty$$

Furthermore, let  $\Pi^{(k_j)}$ ,  $j=1,\ldots,m$ , be independent Poisson processes on  $\mathbf{A}\times[0,1]$  with intensity measure  $d\boldsymbol{\lambda}^{(k_j)}(x,y)=\lambda^{(k_j)}2(k_j-1)y^{2k_j-3}dx\,dy$ . Then

$$(\Xi_n^{(k_1)},\ldots,\Xi_n^{(k_m)}) \stackrel{\mathcal{D}}{\to} (\Pi^{(k_1)},\ldots,\Pi^{(k_m)}).$$

**Proof.** Note that  $\mu(C_n^{(k_j)}) = O(n^{-k_j/(k_j-1)})$  and hence the conditions of Theorem 5.8 are fulfilled. The result is then proved in the same manner as Corollary 5.4.

# 6 Possible extensions, improvements, and applications

#### 6.1 Extensions

A natural extension of this paper is to generalize the results to higher dimensions. All the approximations make use of the probability of covering a number of independently and uniformly distributed particles with some translate of a convex set. In three dimensions this probability can be found in Månsson (1996) and, as shown in Section 3, the probability in an arbitrary dimension can be obtained directly from results in Weil (1990). It should be straightforward, but tedious, to extend the approximation results to an arbitrary dimension.

The starting point can be changed in various directions. An obvious variation is to let the particles on **A** constitute a Poisson process rather than being a fixed number. To derive results, corresponding these of the present paper, in that case should be easy, and it would be surprising if it asymptotically would be any difference. Variations, more difficult to handle, are for instance to let the particles follow some distribution other than the uniform, or to let the sets be non-convex.

#### 6.2 Compound Poisson approximation

The indicators pertaining to k-subsets with common particles are not independent, but have a positive dependence, and the k-subsets which are covered tend to occur in clumps. The more common particles and the larger sets, the stronger is the dependence. Then it seems natural to approximate the process of the positions of the k-aggregates by some process in which clumps are more likely to occur than in the usual Poisson process, and to approximate the number of k-aggregates by some distribution other than Poisson. Natural candidates are the compound Poisson process and the compound Poisson distribution. It seems possible that such approximations can handle larger sets and improve the error bounds, especially in case of large k.

#### 6.3 Statistical applications

As mentioned in the introduction, a common test statistic when testing whether a point pattern originates from a Poisson process is the number of pairs of points closer than some distance r. This number equals W in our terminology, in the case where k=2 and C is a disc with diameter r. An idea is to study if the number of k-subsets,  $k=2,3,\ldots$ , which are covered by a set C, which not necessarily is a disc, would be reasonable as a statistic when testing the randomness of a sample. The crucial element in this matter are the rates of convergence.

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