Correlated Non-Coherent Radar Detection for Gamma-Fluctuating Targets in Compound Clutter

Josef Zuk
Defence Science and Technology Group, Australia

Abstract—This work studies the problem of radar detection of correlated gamma-fluctuating targets in the presence of clutter described by compound models with correlated speckle. If the correlation is not accounted for in a radar model, the required signal-to-interference ratio for a given probability of detection will be incorrect, resulting in over-estimated performance. Although more generally applicable, the is focus on airborne maritime radar systems. Hence K-distributed sea clutter is used as the main example. Detection via square-law non-coherent pulse integration is formulated in a way that accommodates arbitrary partial correlation for both target radar cross-section (RCS) and clutter speckle. The obstacle to including this degree of generality in previous work was the fact that Swerling’s original characterization of the standard RCS fluctuation classes as gamma distributions for the power is not sufficient for the inclusion of both correlation sources (i.e. target and clutter speckle) for gamma-fluctuating targets. An extension of the model is required at the quadrature component (i.e. voltage level), as phase relationships can no longer be neglected. This is addressed in the present work, which not only postulates an extended model, but also demonstrates how to efficiently compute it, with and without a number of simplifying approximation schemes within the framework of the saddle-point technique.

Index Terms—Radar detection, non-coherent, sea clutter, radar cross-section correlation, gamma-fluctuating target, K-distribution.

I. INTRODUCTION

PARAMETRIC modelling is a useful tool for predicting the detection performance of a radar system given a statistical description of the environment. Statistical models of sea clutter are commonly employed in the parametric modelling of detection by maritime surveillance radar [1], [2]. These are often compound models that represent the fast varying speckle component of the clutter with a local Gaussian process which is modulated by a slowly varying texture component. If the radar employs frequency agility, it may be possible to treat the speckle as completely decorrelated from pulse to pulse, however if the radar uses a fixed frequency or the agile frequencies are separated by less than the pulse bandwidth, the speckle correlation must be taken into account.

In this paper, the problem of radar detection using non-coherent pulse integration is considered, where the fluctuations of target RCS are typically modelled in terms of the standard Swerling classes. Varying degrees of temporal correlation of target RCS and clutter speckle fluctuations are relevant considerations for performance modelling of airborne maritime surveillance systems. Therefore, we formulate a modelling scheme that accommodates the simultaneous inclusion of arbitrary partial correlation of both a gamma-fluctuating target RCS and the speckle component of a compound clutter model. In line with the maritime focus of the paper, K-distributed sea clutter is used as the main example.

In the case where the only source of interference is (uncorrelated) thermal noise, general exact calculations for detection probability resulting from non-coherent pulse integration and partially correlated Rayleigh fluctuating target RCS were given by Kanter [3], and later for general gamma-fluctuating targets by Weiner [4]. Various simplified special cases had been previously considered, and a review of these can be found in [5]. A discussion of Kanter’s model in which both the target RCS and clutter returns are present and partially correlated was given in [6]. Subsequently, efficient approximations to the exact but numerically unstable results of Kanter and Weiner have been developed, based on effective number-of-loops concepts [1], [2], and saddle-point techniques [7], [8], [9]. The problem of correlated gamma-fluctuating targets in uncorrelated clutter has also recently been studied in [10].

In previous work on the saddle-point approach to the problem [9], attention was restricted to uncorrelated clutter speckle because the full problem of a correlated gamma-fluctuating target in correlated clutter speckle requires considerations beyond those of computational methodology (saddle-point or otherwise); in particular, an extension of the underlying model must first be established. The classic Swerling models of target RCS fluctuation are specified in terms of the power distributions of the target returns [11]. While this suffices generally for exponentially fluctuating targets, when the target power fluctuations are gamma distributed and when both the target and clutter returns are correlated, the distribution of total returned power depends also on the relative phases of target quadrature components [12] due to a breaking of orthogonal symmetry. A similar issue for clutter is discussed in [13], where it is pointed out that the relative phase between quadrature components is correlated even in the Gaussian problem unless it comprises entirely white noise. A general first-principles model that specifies the joint probability distribution of the quadrature components in a way that reduces correctly to all known special cases and satisfies all expectations has not previously appeared in the literature. A solution to this problem, along with the demonstration of the applicability of the saddle-point technique, is the subject of the present paper. While attention here is focussed on pulse-to-pulse integration, the concepts developed are also applicable to systems employing scan-to-scan integration [14].
The rest of the paper is organized as follows: Section II introduces the problem and the general method of solution using the inverse Laplace transform of the moment generating function. Section III describes the representation of the moment generating function for correlated clutter speckle. Section IV develops the exact physical solution to the problem, designated as a ‘first principles’ model, which is not solvable by analytic means. Section V derives an effective method to evaluate the first-principles model. The resulting effective model serves as a proxy for the first-principles model for computational purposes. A number of approximations are then proposed to further reduce the computational load while maintaining robust performance for all realistic conditions. In Section VI Monte Carlo simulation is employed to demonstrate the high degree of accuracy of the effective method in approximating the first-principles model, in all physically realistic situations. It should be noted that the effective model does not stand on its own: If the first-principles-model were not developed, then there would be no basis upon which to judge whether the effective model is a valid representation of the problem. Section VII then summarizes the key results in the paper.

II. Problem Formulation

To be specific, we consider square-law detection followed by $M$-pulse non-coherent integration, in the presence of multiple partially correlated signal/noise power sources. The test statistic that determines the detection threshold is the summed returned power random variable (RV)

$$Z = \frac{1}{2} \sum_{\eta=1}^{2} \left| \sum_{\alpha} X_{\eta\alpha} \right|^2,$$  

where $X_{\eta\alpha}$ is an $M$-dimensional vector with elements $X_{\eta\alpha m}$, $\eta = 1, 2, \ldots, M$. Vector RVs are indicated by a bold typeface. For $\eta = i, q$, the $X_{\eta\alpha}$ are respectively the in-phase and quadrature components on the $m$-th integrated pulse for signal power source $\alpha$. The source index $\alpha$ will typically range over thermal noise ($n$), target return ($s$) and various types of clutter ($c$). The target RCS fluctuation class indexed by $\nu$ is associated with power source $\alpha$, where $\nu = 1$, in which case $\nu = 2$.

The joint PDF of the vector RV $X_{\eta\alpha}$ is a mean-zero correlated Gaussian with variance $\sigma^2_{\eta\alpha}$ representing the returned power associated with power source $\alpha$, and correlation matrix $C_{\eta\alpha}$. It should be noted that the matrices $C_{\alpha}$ are taken to be correlation matrices, not covariance matrices. Therefore, it is assumed that $\text{Tr} C_{\alpha} = M$. For the thermal noise component, the correlation matrix is the identity: $C_{n} = I$.

It is convenient to introduce the notation $Z_{\text{avg}} \equiv Z/M$ for the pulse-averaged returned power, and $Z_{\text{avg}}^2 \equiv Z_{\text{avg}} / \sigma^2_{\text{avg}}$ for the average power normalized by the mean total interference $\sigma^2_{\text{avg}} = \sigma^2_n + \sigma^2_s$. Also, $S \equiv \sigma^2_s / \sigma^2_n$ will denote the signal-to-interference ratio (SIR), and $q \equiv \sigma^2_s / \sigma^2_n (0 \leq q \leq 1)$ is the clutter-to-interference ratio.

We shall work with the Gauss-Markov correlation model [3], for which the correlation matrices have the symmetric Toeplitz form:

$$C_{\alpha} = \begin{bmatrix} \rho^{|m-n|} & \rho^{|m-n|} & \cdots & \rho^{|m-n|} \\ \rho^{|m-n|} & \rho^{|m-n|} & \cdots & \rho^{|m-n|} \\ \vdots & \vdots & \ddots & \vdots \\ \rho^{|m-n|} & \rho^{|m-n|} & \cdots & \rho^{|m-n|} \end{bmatrix},$$  

where $\rho_n$, $\rho_c$ are the target and clutter correlation coefficients, respectively, such that $0 \leq \rho_n, \rho_c \leq 1$. Nevertheless, the methodology that we develop works equally with any other Toeplitz correlation structure.

By considering the various known special cases, such as correlated exponentially fluctuating target power ($\kappa = 1$) with correlated clutter speckle [8], [12], and correlated gamma-fluctuating target power ($\kappa > 1$) with uncorrelated clutter speckle [9], one can infer that the mean and variance of the total returned power in the general case should be given, respectively, by

$$\langle Z_{\text{avg}}^2 / \sigma^2_{\text{avg}} \rangle = 1 + S,$$  

and

$$\text{var} \left( Z_{\text{avg}}^2 / \sigma^2_{\text{avg}} \right) = \frac{1 - q^2}{M} + \frac{q^2}{L} + S^2 \left( \frac{1 - q}{M} + \frac{q}{N} \right),$$  

where

$$L \equiv M^2 / \text{Tr} \{ C_n^2 \}, \quad B \equiv M^2 / \text{Tr} \{ C_c^2 \}, \quad N \equiv M^2 / \text{Tr} \{ C_n C_c \}. $$

The quantity $\zeta$ represents the effect of compound clutter texture, and is given by

$$\zeta = 1 + (L + 1) \text{var} (U) = 1 + (L + 1) / \nu,$$  

with the second equality holding for K-distributed compound clutter, where the texture RV $U$ is described by a unit-mean gamma distribution with shape parameter equal to the clutter texture shape parameter $\nu$, in which case $\text{var} (U) = 1/\nu$. Standard Rayleigh clutter corresponds to $\zeta = 1$, obtained in the limit $\nu \to \infty$. It should be noted that the considerations in this work remain valid for any compound clutter model. In the general case, $U$ is the random variable that realizes the clutter texture distribution in any compound clutter model.

The simplest way to understand [7] is to set $S = 0$ and $q = 1$. Then $Z_{\text{avg}}^2 / \sigma^2_{\text{avg}} = UV$, where the clutter speckle RV $V$ is given by the correlated sum of $M$ unit-mean exponential RVs and scaled by $M$). Since the second moment of the compound
RV factorizes into texture and speckle components, we have \( \text{var}(Z_{\text{ng}}/\alpha^2) = m_2(U) m_2(V) - 1 \), given that \( \langle U \rangle = \langle V \rangle = 1 \). The result follows upon observing that \( \text{var}(V) = 1/L \) [3, Eq. 65].

The challenge is to derive a general model that reproduces this mean and variance. It is worth mentioning at this point that the related problem of including partial correlation of clutter speckle in cases where a non-Rayleigh overall clutter amplitude is being modelled was successfully addressed by the introduction of the spherically invariant process (SIRP) [13], of which the compound clutter paradigm is a special case. However, this approach is not applicable to the correlation of non-Rayleigh target amplitudes because introducing a SIRP leads to an increase in power variance, as can be seen for the clutter contribution in (5) and (7) by comparing a finite shape parameter \( \nu \) with \( \nu \rightarrow \infty \); On the other hand, moving the RCS fluctuation class index \( \kappa \) to values greater than unity will necessarily decrease the variance.

Compound clutter is incorporated in the usual way [8, 12] by promoting the mean clutter power to a random variable according to the mapping \( \alpha^2 \rightarrow \bar{\alpha}^2 U \) where the RV \( U \) generates the clutter texture distribution with unit mean. The overall mean interference power becomes

\[
\sigma_i^2 = \sigma_n^2 + \langle \alpha^2(U) \rangle_U = \sigma_n^2 + \bar{\alpha}^2 .
\] (8)

The texture expectation is best left to the end, after recovering the survival function from the MGF. Thus, if \( F_{cc}(v; S) \) denotes the survival function for the total returned interference power \( v \) in the presence of both a target and compound clutter, and we also use the notation \( F_{spk}(v; S, q) \), where \( q \equiv \bar{\alpha}^2 / \sigma_n^2 \) is the clutter-to-interference ratio, to denote the survival function incorporating clutter speckle but without the texture, then \( F_{cc}(v; S) \) is given by

\[
F_{cc}(v; S) = \left( F_{spk}(v; S, qU) \right)_U = \int_0^\infty du f_{\nu}(u) F_{spk}(v; Su) ,
\] (9)

where \( f_\nu(u) \) denotes the clutter texture PDF. For K-distributed clutter, this is a unit-mean gamma distribution with shape parameter \( \nu \) [1]. The texture integration is most efficiently performed by means of Gaussian quadrature.

The probability of detection of a target on a single scan \( P_{d}(S) \) when its SIR is known to be \( S \), is given by the value of the survival function (also known as the complementary cumulative distribution function) for the total returned signal plus interference power at the selected detection threshold \( v = v_b \). The threshold is determined as the point at which the survival function for the total returned interference power assumes the value that corresponds to a pre-determined desired probability of false alarm \( P_{\text{FA}} \). Thus, we eliminate \( v_b \) in the following pair of equations:

\[
P_{d}(S) = F_{cc}(v_b; S) , \quad P_{\text{FA}} = F_{cc}(v_b; 0) .
\] (10)

In the saddle-point method, just as for the PDF, one computes the survival function in terms of the inverse Laplace transform of the MGF for its underlying distribution, by integrating along a suitably constructed contour in the complex plane [7].

Previous work [9] applied saddle-point techniques to the problem of correlated targets in uncorrelated clutter speckle, using an existing model for the problem. The present work, in part, also applies saddle-point techniques, but to a model that must first be developed, as there is no extant model that addresses the problem of correlated gamma-fluctuating target in correlated clutter speckle.

III. CORRELATED CLUTTER

Weiner’s model [4] is applicable to correlated gamma-fluctuating targets embedded in uncorrelated clutter speckle. In this section, we generalize this by establishing the representation of the MGF that incorporates speckle correlation, and discuss the complications encountered when trying to evaluate it according to the usual specification of the Swerling RCS models.

We assume that there are three sources of returned power \( \alpha = n, c, s \), corresponding to thermal noise, surface clutter and target RCS, respectively, and that the correlated gamma-fluctuating target is embedded in correlated clutter (belonging to the \( \kappa = 1 \) speckle fluctuation class). The \( M \)-dimensional vector RV \( X_v \), composed of one element for each of the \( m = 1, 2, \ldots, M \) pulses, will denote the in-phase quadrature component for the associated power sources.

The MGF \( M_Z(s) \) of the returned-power RV \( Z \) is given by

\[
\sqrt{M_Z(s)} = \left( e^{\frac{s}{2} (\sum_\alpha X_\alpha)} \right)_X = \int_{-\infty}^{+\infty} \frac{d\alpha}{(2\pi s)^{M/2}} e^{-\alpha \cdot \alpha/(2s)} \prod_{\alpha = n, c, s} (e^{-iU X_\alpha})_X ,
\] (11)

where we have split off the non-Gaussian target component \( X_s \), evaluated the expectations over the clutter and noise components, and set

\[
Q(s) \equiv (1 + \sigma_n^2) I + \sigma_c^2 C_c .
\] (12)

The square-root on the LHS of the equation arises from the square on the RHS in (12).

Now, since \( Q = R^T A R \), we may write \( A = \text{diag}(\lambda_1, \ldots, \lambda_M) \) and \( R^T R = I \). Then we have, for any vector \( J \), the representation

\[
\int_{-\infty}^{+\infty} \frac{d\alpha}{(2\pi s)^{M/2}} e^{-\alpha^T Q \alpha/2, e^{-iu^T J}} = \int_{-\infty}^{+\infty} \frac{d\alpha}{(2\pi s)^{M/2}} e^{-u^T \Lambda u/2, e^{-iu^T R J}} = \frac{1}{\sqrt{\text{det} Q}} \exp \left[ -\sum_{m=1}^{M} (RJ)^2_m / (2\lambda_m) \right] ,
\] (13)
where the first equality is due to a rotational change of integration variables \(u' = Ru\), and we note that

\[
\sum_{m=1}^{M} (RJ_m^2)/\lambda_m = \sum_{m,a,b=1} J_a (R_{am}^\dagger \lambda_m R_{mb}) J_b
\]

\[
= J^T Q^{-1} J ,
\]

which leads to the result

\[
\sqrt{M_Z(s)} = \frac{1}{s^{M/2}/\det Q(s)} \left< e^{-\frac{1}{2}X_s^T Q^{-1}(s)X_s} \right>_{X_s},
\]

upon making the identification \(J \equiv X_s\).

It is convenient to introduce a scaled target RV \(X_s\), according to \(X_s = \sigma_s X_s\), so that each power component \(X_{sm}^2\) has unit mean. For RCS fluctuation class \(\kappa\), each \(X_{sm}^2\) has a marginal distribution given by the unit-mean gamma distribution with shape parameter \(\kappa/2\), i.e., \(X_{sm}^2 \sim \Gamma(\kappa/2, 2/\kappa)\) for all \(m = 1, 2, \ldots, M\) \([11]\).

For uncorrelated clutter speckle, \(Q\) becomes proportional to the unit matrix \(Q = \xi^2 I\), with

\[
\xi^2 \equiv 1/s + \sigma_\alpha^2 + \sigma_\beta^2 = 1/s + \sigma_\alpha^2.
\]

This yields the representation

\[
\sqrt{M_Z(s)} = \frac{1}{(s \xi^2)^{M/2}} \left< \exp \left( \frac{\sigma_\alpha^2}{2s \xi^2} X_s^2 \right) \right>_{X_s},
\]

which reproduces the problem considered in the previous work \([9]\) when combined with the expression \([1]\)

\[
\left< \exp \left( -\frac{\sigma_\alpha^2}{2s \xi^2} X_s^2 \right) \right>_{X_s} = \left[ \det \left( I + \frac{\sigma_\alpha^2}{\kappa \xi^2} C_s \right) \right]^{-\kappa/2}. \tag{18}
\]

The higher-order Swerling models \((\kappa > 1)\) are defined by giving the common distribution of the squares of the quadrature components \(X_{sm}^2\) \([11]\), but the signs are left unspecified. In the presence of partial correlation, however, the signs of the target quadrature components do not drop out and, therefore, the specification is incomplete. Equivalently, one may note that while \(X_s^2\) is rotationally invariant, the quadratic form \(X_s^T Q^{-1} X_s\) is not.

IV. A FIRST-PRINCIPLES MODEL

In this section, we review the derivation of Weiner’s model for gamma-fluctuating targets in uncorrelated interference, and use it to motivate a generalization to accommodate the presence of clutter with an arbitrarily correlated speckle component. The aim is to derive a first-principles model, i.e., one whose degrees of freedom correspond to basic observable physical quantities. In the present context, these are provided by the electromagnetic field impinging upon the antenna aperture or, equivalently, by the voltage variations induced because of it. Due to the stochastic nature of the detection problem, the relevant physics inputs to the model are the probability distributions of the two orthogonal quadrature components.

According to Weiner’s formulation \([3]\), in fluctuation class \(\kappa\), the scaled target power RV may be characterized as

\[
X_s^2 = \frac{1}{\kappa} \sum_{k=1}^{\kappa} H_k^\dagger H_k,
\]

where, for each \(k = 1, 2, \ldots, \kappa\), the vector RV \(H_k\) is a zero-mean unit-variance Gaussian with common correlation matrix \(C_s\). The associated probability measure \(F_{H}(dh) = P_{H}(h)dh\) is thus given by

\[
F_{H}(dh) = \frac{dh}{[(2\pi)^M \det C_s]^{\kappa/2}} \exp \left\{ -\frac{1}{2} \sum_{k=1}^{\kappa} h_k^T C_s^{-1} h_k \right\}, \tag{20}
\]

where \(dh\) is the product measure in \(\kappa M\) dimensions

\[
dh = \prod_{k=1}^{\kappa} \prod_{m=1}^{M} dh_{k,m}, \tag{21}
\]

and where \(h_{k,m}\) denotes the \(m\)-th component of the vector \(H_k\). We observe the explicit orthogonal invariance \(F(d(Rh); RC_sR^T) = F(dh; C_s)\), where \(F(dh; C_s) \equiv F_{H}(dh)\) and \(Rh \equiv \{Rh_1, Rh_2, \ldots, Rh_k\}\). If we write \(T \equiv X_s^2\), then the induced probability measure \(F_T(dt) = P_T(t)dt\) on the space of functions of the scalar RV \(T\) is given by

\[
F_T(dt) = \int_{R^M} dh P_{H}(h) \delta(t - \frac{1}{\kappa} \sum_{k=1}^{\kappa} h_k^T h_k) dt , \tag{22}
\]

where \(\delta(t)\) is the Dirac delta function. Appealing to the Fourier transform of the delta function

\[
\delta(t) = \int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} e^{i\omega t} , \tag{23}
\]

it follows immediately that

\[
P_T(t) = \int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} e^{i\omega t} \left[ \det \left( I - i\omega \frac{2\pi}{\kappa} C_s \right) \right]^{-\kappa/2} , \tag{24}
\]

which depends only on the eigenvalues of \(C_s\), as expected due to the invariance under the mapping \(C_s \mapsto RC_s R^T\). When \(C_s = I\), we recognize, in the integrand, the characteristic function for the distribution \(T \sim \Gamma(M\kappa/2, 2/\kappa)\), which corresponds to shape parameter \(M\kappa/2\) and mean \(M\). And if we write

\[
T_m = \frac{1}{\kappa} \sum_{k=1}^{\kappa} H_{mk}^2 , \tag{25}
\]

so that \(T = \sum_{m=1}^{M} T_m\), then \(T_m \sim \Gamma(\kappa/2, 2/\kappa)\) for each \(m = 1, 2, \ldots, M\). The relevant question now is how to extend this discussion in order to construct a compatible probability measure on the space of functions of the vector RV \(X_s\).

At this point, we can see that \([18]\) may be established by appealing to the orthogonal invariance of the expectation on the LHS. Specifically, the expectation

\[
\left< \exp \left( -\frac{\sigma_\alpha^2}{2s \xi^2} X_s^2 \right) \right>_{X_s} = \left< \exp \left( -\frac{\sigma_\alpha^2}{2s \xi^2} \sum_{k=1}^{\kappa} H_k^T H_k \right) \right>_{H} \tag{26}
\]

is unchanged under the mapping \(H_k \mapsto RH_k\) for any orthogonal matrix \(R\), provided that it is accompanied by the mapping \(C_s \mapsto RC_s R^T\). To take advantage of this, we write the target correlation matrix in diagonalized form as \(C_s = R_s^T \Lambda_s R_s\), where \(\Lambda_s \equiv \text{diag}(\gamma_1, \gamma_2, \ldots, \gamma_M)\) and \(R_s^T R_s = I\), and introduce the rotated vector RVs \(J_k \equiv R_s H_k\) whose elements are independent zero-mean Gaussians such that the \(J_{mk}\) have
common variance $\gamma_m^k$ for $k = 1, 2, \ldots, \kappa$. It follows that the sums
\[ V_m \equiv \frac{1}{\kappa} \sum_{k=1}^{\kappa} J_{mk}^2 \] (27)
are independent and gamma-distributed with respective means $\gamma_m^k$. Then, we explicitly have
\[ \exp \left( -\frac{\sigma^2}{2\kappa} \sum_{m=1}^{M} V_m \right) = \prod_{m=1}^{M} \left( 1 + \frac{\sigma^2}{\kappa \xi \gamma_m^k} \right)^{-\kappa/2} \], (28)
and this is equivalent to (18).

We may apply the same rotational change of variable scheme to the problem with correlated clutter, where $Q$ is a general matrix, provided we can assume that the, yet unspecified, mapping $(H_1, H_2, \ldots, H_s) \mapsto \hat{X}_s$ satisfies $(RH_1, RH_2, \ldots, RH_s) \mapsto R\hat{X}_s$ for any orthogonal matrix $R$. This requirement is clearly consistent with (15) but its existence for $\kappa > 1$ is not clear. In this case, the argument of the expectation in (15) is not rotationally invariant, but the induced probability measure on $\hat{X}_s$ will be invariant under the simultaneous mappings $X_s \mapsto R_s\hat{X}_s$, $C_s \mapsto R_sC_sR_s^T$. Thus, setting $Y \equiv R_s\hat{X}_s$, we obtain
\[ \mathbb{E} \left[ \exp \left( -\frac{\sigma^2}{2} \hat{Y}^T Q^{-1} \hat{Y} \right) \right] = \mathbb{E} \left[ \exp \left( -\frac{\sigma^2}{2} \hat{Y}^T Q^{-1} \hat{X}_s \right) \right] = \exp \left( -\frac{\sigma^2}{2} \hat{Y}^T (R_sQ^{-1}R_s^T) \hat{Y} \right) = \exp \left( -\frac{\sigma^2}{2} \hat{Y}^T (L_sQ^{-1}L_s^T) \hat{Y} \right), \] (29)
where $Y \equiv \sqrt{\lambda_s} \hat{Y}$, and $L_s \equiv \sqrt{\lambda_s} R_s$ so that $C_s = L_s^T L_s$. We note that elements of the vector RV $Y$ are independent and identically distributed (iid) such that $Y_m^2 \sim \Gamma(\kappa/2, 2/\kappa)$ for all $m = 1, 2, \ldots, M$.

The problem of an uncorrelated gamma-fluctuating target embedded in correlated clutter is obtained when $L_s = I$. Therefore, we see that introducing target correlation in the presence of correlated clutter simply amounts to making the mapping $Q^{-1} \mapsto L_sQ^{-1}L_s^T$ in the calculation of the target expectation. When $C_s = I$, the only sensible option is for each of iid RVs $\hat{X}_m$ to be described by a two-sided Nakagami distribution whose PDF is given by
\[ f(x; \nu, \Omega)dx = \frac{1}{\Gamma(\nu)} e^{-\left( x\sqrt{\nu}/\Omega \right)^2} \left[ x\sqrt{\nu}/\Omega \right]^{2\nu-1} d(x\sqrt{\nu}/\Omega), \] (30)
with fading parameter $\nu = \kappa/2$ and scale parameter $\Omega = 1$, and this becomes Gaussian for $\kappa = 1$. The characteristic function is given by [15]
\[ \hat{f}(\xi; \nu, \Omega) \equiv e^{-iu\hat{X}} = \frac{\Gamma(2\nu)}{2\nu \Gamma(\nu)} \exp \left( -\frac{\Omega}{8\nu} u^2 \right) \times \left[ D_{-2\nu} \left( iu \sqrt{\frac{\Omega}{2\nu}} \right) + D_{-2\nu} \left( -iu \sqrt{\frac{\Omega}{2\nu}} \right) \right], \] (31)
where $D_{-\nu}(z)$ is a parabolic cylinder function [16]. Regardless of whether it is possible to find a suitably rotationally covariant mapping $(H_k) \mapsto \hat{X}_s$, we shall adopt (29) as the definition of a candidate first-principles model: that is, we set $X_s \equiv L_s^T \hat{Y}$, and furthermore postulate that, even for a general target correlation matrix $C_s$, the components of the vector RV $Y$ are iid and generated by the two-sided Nakagami distribution.

A. Statistical Properties of the First-Principles Model

The aim of this section is to demonstrate that the model proposed in the foregoing section yields the required mean and variance as given by (4)–(6). Recalling that the mean and variance are respectively given by the first and second derivatives of the cumulant generating function (CGF) evaluated at the origin, we proceed by writing down the CGF $K(s)$ for the model and making a Taylor expansion about $s = 0$ up to second order.

Thus, in order to compute the mean and variance arising from (15), let us introduce the functions
\[ K_0(s) = -\ln \det(s\cdot Q(s)), \]
\[ K_1(s) = 2 \ln \left( \exp \left( -\frac{\sigma^2}{2} \hat{X}^T Q^{-1}(s) \hat{X} \right) \right) \] (32)

Then, the cumulant generating function for the model defined by (15) is given by
\[ K(s) = \ln \langle e^{-s\hat{Z}_{ns}} \rangle_{\hat{Z}_{ns}} = K_0(s) + K_1(s). \] (33)

The function $K_0(s)$ generates the contribution to the CGF that is independent of the SIR. Thus, performing a Taylor expansion in $s$,
\[ K_0(s) = -s(\sigma_n^2 + \sigma_c^2) + s^2 \left( \frac{\sigma_n^4}{M} + \frac{\sigma_c^4}{M^2} \right) \text{Tr}(C_n^2) + \frac{2\sigma_n^2\sigma_c^2}{M} + O(s^3). \] (34)

Also performing a Taylor expansion in $s$ for the SIR-dependent term yields
\[ K_1(s) = -s\sigma_n^2 \left( \hat{X}^T Q^{-1}(s) \hat{X} \right) \ ] \ X_s \]
\[ + \frac{\sigma_n^4}{4} \left[ \left( \hat{X}^T Q^{-1}(s) \hat{X} \right)^2 \right] \ ] \ X_s \]
\[ + s^2 \left( \frac{\sigma_n^4}{M} + \frac{\sigma_c^4}{M^2} \right) \text{Tr}(C_n^2) + \frac{\sigma_n^2\sigma_c^2}{M} + O(s^3), \]
\[ + \frac{\sigma_n^2\sigma_c^2}{M} + O(s^3), \] (35)
and we note that
\[ Q^{-1}(s) = \frac{1}{1 + s(\sigma_n^2 + \sigma_c^2)} = sI - s^2(\sigma_n^2I + \sigma_c^2C_n) + O(s^3), \] (36)
which yields
\[ K_1(s) = -s\sigma_n^2 \left( \hat{X}^T Q^{-1}(s) \hat{X} \right) \ ] \ X_s \]
\[ + s^2 \left( \frac{\sigma_n^4}{M} + \frac{\sigma_c^4}{M^2} \right) \text{Tr}(C_n^2) + \frac{\sigma_n^2\sigma_c^2}{M} + O(s^3). \] (37)

A matrix in the denominator, as in (36), denotes the inverse.
By construction, \( n \) [29] holds in distribution for the first-principles model. This is easily confirmed, based on the observation that
\[
H \sim \mathcal{N}(0,1) \Rightarrow H^2 \sim \Gamma(1/2,2) .
\] (38)
It follows immediately that
\[
\langle \hat{X}_m^T \hat{X}_n \rangle_{\hat{X}_m} = \text{Tr}\{C_s\}, \quad \text{var}(\hat{X}_m^T \hat{X}_n) = \frac{2}{\kappa} \text{Tr}\{C_s^2\} .
\] (39)
From this, we see that the mean and variance are given as in (4)–(6), but with
\[
N = M^2 / \langle \hat{X}_m^T C_s \hat{X}_n \rangle_{\hat{X}_m} .
\] (40)
To compute this expectation, we recall that the RVs \( \hat{Y}_m \), \( m = 1, 2, \ldots, M \) are iid such that \( \hat{Y}_m^2 \sim \Gamma(\kappa/2, 2/\kappa) \) and with random signs (i.e. randomly positive or negative, so that \( \text{sgn}(\hat{Y}_m) = \pm 1 \) is by construction a Rademacher RV), and set \( \hat{X}_m \equiv L_m^T \hat{Y} \), \( C_s = L_m^T L_m \). Accordingly,
\[
\langle \hat{X}_m^T C_s \hat{X}_n \rangle_{\hat{X}_m} = \sum_{a,b=1}^M [L_s]_{am} [L_s]_{bn} \langle Y_a Y_b \rangle_{\hat{Y}} \nonumber
\]
\[
= \sum_{a,b=1}^M [L_s]_{am} [L_s]_{bn} \delta_{ab} \nonumber
\]
\[
= [L_s^T L_s]_{mn} \nonumber
\]
\[
= [C_s]_{mn} .
\] (41)
This means that we obtain \( \langle \hat{X}_m^T C_s \hat{X}_n \rangle_{\hat{X}_m} = \text{Tr}\{C_s C_s\} \) as desired. The upshot of all this is that the candidate first-principles model reproduces the expected mean and variance for the returned power, and \( \langle f \rangle \) holds, but the rotational covariance assumption that initially led to it is redundant.

B. Model Summary

In summary, the first-principles model is defined by specifying that the (normalized) target quadrature component is given by \( X_s \equiv L_s^T Y \), where the iid vector RV \( Y \) has elements distributed according to the two-sided Nakagami distribution of (30), and the matrix \( L_s \) is obtained from the eigenvalue decomposition of the target correlation matrix according to
\[
C_s = R_s^T \Lambda_s R_s , \quad L_s \equiv \sqrt{\Lambda_s} R_s .
\] (42)
Hence, the MGF of the first-principles model is given by
\[
\mathcal{M}_Z(s) = \frac{1}{s^M \det Q(s)} \left( \exp \left( -\frac{\sigma_n^2}{2} Y^T (L_s Q^{-1}(s) L_s^T) Y \right) \right)^2 .
\] (43)
In the Appendix, we discuss some minor shortcomings of the first-principles model. We also show that it may be computed exactly for the special case of a fully correlated target.

Compound clutter may be incorporated as described at the end of Section II by making the \( Q \)-matrix of (12) depend on the clutter texture integration variable \( u \), according to
\[
Q(s; u) \equiv (1/s + \sigma_n^2) I + u \sigma_c^2 C_s ,
\] (44)
or, equivalently, the normalized form
\[
Q(s; u)/\sigma_n^2 = \left( \frac{1}{s \sigma_n^2} + 1 - q \right) I + qu C_s .
\] (45)
The calculations proceed for an arbitrary value of \( u \), until the texture integration is performed at the end, as indicated in (9).

V. Effective Model

A ‘first-principles’ model for the problem of a correlated gamma-fluctuating target in correlated clutter would be characterized by the specification of the joint probability distribution for the quadrature RVs \( \hat{X}_s \) that appear in (15) such that \( \langle f \rangle \) holds. However, having such a joint probability distribution does not necessarily, or even likely, lead to a convenient closed-form MGF to which saddle-point techniques could be easily applied. This is the case with the proposal of the preceding section, as defined by (43), which appears to be analytically intractable. So in this section, we propose an alternative ‘effective’ model — one whose outputs are expected to be that same as those of the first-principles model, though not specified in terms of physical degrees of freedom — that is amenable to analytical treatment. Specifically, our approach is to postulate an explicit MGF that reduces to the expected results in all known computable special cases. This MGF would then define an ‘effective model’ for the system. We examine its relationship with the previously defined first-principles model, show how to compute the survival function from the MGF using saddle-point techniques and, finally, present two approximation schemes that reduce computational load significantly while maintaining a high degree of accuracy.

Let us consider the following ‘effective’ MGF for the problem with correlated clutter:
\[
\mathcal{M}_{Z_{avg}}(s) = \frac{\left[ \det (I + (s/M) (\sigma_n^2 + \sigma_c^2 C_s)) \right]^{\kappa-1}}{\left[ \det (I + (s/M) (\sigma_n^2 + \sigma_c^2 C_s + \sigma_c^2 C_s/k)) \right]^\kappa},
\] (46)
which correctly reproduces the special cases (i) \( \sigma_n^2 = 0 \), (ii) \( \sigma_c^2 = 0 \), (iii) \( \rho_c = 0 \), and (iv) \( \kappa = 1 \). It also satisfies the essential requirement of yielding the expected mean and variance, as given by (4)–(6). To see that, we construct the CGF
\[
\mathcal{K}(t) = \ln \mathcal{M}_{Z_{avg}}(t) = \ln \langle e^{-tZ_{avg}} \rangle_{Z_{avg}},
\] (47)
which, for the present problem, reads
\[
\mathcal{K}(t) = (\kappa - 1) \text{Tr} \left\{ \ln \left[ I + (t/M) (\sigma_n^2 + \sigma_c^2 C_s) \right] \right\}
\]
\[
- \kappa \text{Tr} \left\{ \ln \left[ I + (t/M) (\sigma_n^2 + \sigma_c^2 C_s + \sigma_c^2 C_s/k) \right] \right\}
\]
\[
= -t \left( \sigma_n^2 + \sigma_c^2 + \sigma_n^2 \right) + \frac{t^2}{2} \left[ \frac{\sigma_n^2}{M} + \frac{\sigma_c^2}{M^2} \text{Tr}\{C_s\} + \frac{\sigma_c^4}{\kappa M^2} \text{Tr}\{C_s\} \right]
\]
\[
+ 2 \left( \frac{\sigma_n^2 \sigma_c^2}{M} + \frac{\sigma_n^2 \sigma_c^2}{M} + \frac{\sigma_n^4}{M} \text{Tr}\{C_s\} \right) + O(t^3),
\] (48)
and recall that the mean and variance are given by the first and second derivatives of the CGF evaluated at the origin. This expression leads directly to the results given in (4) and (5) for the returned power mean and variance, respectively, noting that we must set \( \nu = \infty \) for the clutter shape parameter as we have
not yet introduced the compound clutter texture. The form of the MGF in [46] is an ansatz (i.e., trial form) motivated by previous work on the heuristic generalized Dalle Mese Giulii (DMG) approximation [12].

In order to make contact with the previous approach, we can write

\[ M_Z(s) = \frac{1}{s^M \det Q(s)} \left[ \frac{\det Q(s)}{\det (Q(s) + \sigma^2 C/M/k)} \right]^\kappa \]

\[ = \frac{1}{s^M \det Q(s)} \left[ \frac{\det \left( I + \frac{\sigma^2}{k} L_s Q^{-1}(s)L_s^T \right)}{\det \left( I + \frac{\sigma^2}{k} L_s Q^{-1}(s)L_s^T \right)} \right]^{-\kappa}. \]  

(49)

We see that reconciling the first-principles approach of [15] and [29] with the effective model of [46] would require that

\[ \left\langle \exp \left( -\frac{1}{2} \hat{Y}^T A \hat{Y} \right) \right\rangle_Y = \left[ \det (I + A/k) \right]^{-\kappa/2}, \]

(50)

for any positive semi-definite symmetric matrix A. We should note that this statement does indeed hold for any such diagonal matrix A, or when \( \kappa = 1 \) in which case \( \hat{Y} \) is a Gaussian RV. However, this relationship will not hold in general: The RHS manifestly only depends on the eigenvalues of A. The LHS will depend only on the eigenvalues if the probability measure on \( \hat{Y} \) is rotationally invariant, which is not expected to be the case that \( \kappa > 1 \). Nevertheless, it is useful to consider the special case where, as is true in the generalized DMG model [12] (discussed in more detail in a later section), the matrices \( C_c \) and \( C_t \) commute, and are thus simultaneously diagonalizable. Then, upon writing \( C_c = R^{-1}A_cR \) and \( C_t = R^{-1}A_tR \) so that \( L_s = \sqrt{\kappa}R \), we have

\[ A = \sigma^2 L_s Q^{-1}(s)L_s^T = \frac{\sigma^2 A_c}{1/s + \sigma^2 A_c + \sigma^2 A_t}, \]  

(51)

which is diagonal, and so the equivalence holds. It follows that the DMG model is simultaneously an approximation to the effective model and any valid first-principles model.

It is also true, in the general case, that the correlation matrices \( C_c \) and \( C_t \) commute asymptotically in the limit as \( M \to \infty \), in the sense that the weak matrix norm of the commutator vanishes in this limit, provided that they are symmetric Toeplitz and satisfy the summability conditions [17]

\[ \sum_{m=1}^\infty ||C_{c,m}| < \infty, \sum_{m=1}^\infty ||C_{t,m}| < \infty. \]

(52)

Thus, (51) holds asymptotically, from which it follows that the effective model is (at least) asymptotically equivalent to the first-principles model. Finally, we remark that, when \( M = 2 \), all correlation matrices will have the circulant form

\[ C = \begin{pmatrix} 1 & \rho \\ \rho & 1 \end{pmatrix}, \]

(53)

and therefore all such matrices commute. We conclude that the first-principles and effective models must be identical for \( M = 2 \).

A. Saddle-point Method for the Effective Model

A robust and accurate way of computing the survival function for a distribution from its MGF is provided by saddle-point techniques, which are especially effective when the MGF is a rational function. This approach has previously been developed for the case of a correlated gamma-fluctuating target embedded in uncorrelated clutter speckle [9], and is easily extended to the general problem.

In the presence of clutter texture, the speckle MGF for the effective model can be expressed as

\[ Z_{\text{avg}}^{spk}(s; u) = \frac{[\det (I + (s/M)(1 - q + qu C_c))|^{\kappa - 1}}{[\det (I + (s/M)(1 - q + qu C_c + SC_s/k))]^s}. \]  

(54)

where \( u \) denotes the normalized texture integration variable. It is convenient to introduce an aggregated target/clutter correlation matrix according to

\[ C_{\text{avg}}(u) \equiv \frac{1}{qu S/k} [qu C_c + (S/k) C_s] \],

(55)

such that \( \text{Tr} C_{\text{avg}} = M \). Then, we can write

\[ Z_{\text{avg}}^{spk}(s; u) = \frac{\prod_{m=1}^M [1 + a_{qm}(u)s]^{\kappa - 1}}{\prod_{m=1}^M [1 + a_m(u)s]^s}, \]

(56)

with

\[ a_{qm}(u) = \left[ 1 - q + qu \gamma^c_m \right] / M, \]

(57)

\[ a_m(u) = \left[ 1 - q + (qu + S/k) \gamma^e_m(u) \right] / M, \]

where

\[ \{ \gamma^c_m: m = 1, 2, \ldots, M \} = \text{eig}(C_c), \]

(58)

\[ \{ \gamma^e_m(u): m = 1, 2, \ldots, M \} = \text{eig}(C_{\text{avg}}(u)). \]

It is useful to note that

\[ \sum_{m=1}^M (a_m - a_{qm}) = S / \kappa. \]

(59)

The saddle-point equation, for the saddle-point (SP) \( s = s_0 \), is obtained as the extremum \( \Phi'(s) = 0 \) of Helstrom’s phase function [7]

\[ \Phi(s; u) \equiv \ln Z_{\text{avg}}^{spk}(s; u) - \ln(-s + su), \]

(60)

and reads

\[ v = \frac{1}{s_0} + \kappa \sum_{m=1}^M \frac{a_m(u)}{1 + a_m(u)s_0} - \frac{a_{qm}(u)}{1 + a_{qm}(u)s_0}. \]

(61)

The integration contour for the inverse Laplace transform of the MGF that yields the survival function is taken to be the steepest descent path (SDP), which is defined by \( \ln \Phi(s) = 0 \) and passes through the saddle-point \( s_0 \). Next, for convenience, we set \( a_0 \equiv a_{q0} \equiv +\infty \) and define

\[ c_m \equiv \frac{1}{s_0 v + v/a_m(u)}, \quad c_{qm} \equiv \frac{1}{s_0 v + v/a_{qm}(u)}, \]

for \( m = 0, 1, \ldots, M \). Then, with \( z \equiv (s_0 - s)v \), we have the associated \( \tau \) phase function [9] defined by

\[ \tau(z) \equiv \Phi(s_0; v) - \Phi(s_0 - z/v; v), \]

(63)
and evaluated as
\[ \tau(z) = z + \kappa \sum_{m=0}^{M} \ln(1 - c_m z) - (\kappa - 1) \sum_{m=0}^{M} \ln(1 - c_{qm} z) \]
\[ = - \sum_{n=2}^{\infty} \frac{r_n}{n} z^n, \tag{64} \]
with coefficients
\[ r_n = \sum_{m=0}^{M} \left[ \kappa c_m^n - (\kappa - 1) c_{qm}^n \right], \tag{65} \]
for \( n = 1, 2, \ldots \), arising in its Taylor expansion. Inversion of this equation (to obtain \( z(\tau) \)), followed by integration over \( \tau \) leads to the exact representation of the speckle survival function
\[ F_{spk}(v; qu) = F_{sp}(v; qu) \sqrt{\frac{2\gamma}{\pi}} \int_{0}^{\infty} d\tau e^{-\tau} \Im z(\tau), \tag{66} \]
assuming that \( s_0 < 1 \) with the factor
\[ F_{sp}(v; qu) = \frac{1}{\sqrt{2\gamma \pi}} e^{\phi(s_0, v, u)} \tag{67} \]
constituting the basic saddle-point approximation. It only remains to perform the clutter texture integration to obtain the full survival function
\[ F_{cc}(v) \simeq \sum_{\ell=1}^{L} w_{\ell} F_{spk}(v; qu_{\ell}), \tag{68} \]
where the \( w_{\ell}, u_{\ell} \) are the weights and nodes, respectively, of an appropriate Gaussian quadrature of order \( L \).

The Padé-adapted version of the \( \tau \)-phase function reads
\[ \tau(z) = z + \ln(1 - c_0 z) \]
\[ + \kappa \left[ (M - 1) \ln(1 - \bar{c} z) + \ln(1 - c_M z) \right] \]
\[ - (\kappa - 1) \left[ (M - 1) \ln(1 - \bar{c} q z) + \ln(1 - c_{qM} z) \right] \]
\[ - \sum_{n=2}^{\infty} \bar{r}_n z^n, \tag{69} \]
with
\[ \bar{c} \equiv \frac{1}{M-1} \sum_{m=1}^{M-1} c_m, \quad \bar{c}_q \equiv \frac{1}{M-1} \sum_{m=1}^{M-1} c_{qm}, \tag{70} \]
and
\[ \bar{r}_n \equiv \sum_{m=1}^{M-1} \left[ \kappa (c_m^n - \bar{c}^n) - (\kappa - 1) (c_{qm}^n - \bar{c}_q^n) \right]. \tag{71} \]
In the special case of uncorrelated clutter speckle (i.e. the problem previously studied in [9]), we have \( c_{qm} = \bar{c} q \) for all \( m = 1, 2, \ldots , M \). In this case, (69) reduces to equation (50) of [9]. Also, the SP equation implies that \( r_1 = 1 \), which expands to the relationship
\[ c_0 + \kappa \left[ c_m + (M - 1) \bar{c} \right] - (\kappa - 1) \left[ c_{qm} + (M - 1) \bar{c}_q \right] = 1. \tag{72} \]

Using this, we can confirm that the contribution to the RHS of (69) linear in \( z \) vanishes, as expected.

The Padé approximation comprises approximating the infinite series in (69) by a low order Padé approximant, as described in [9]. This serves to greatly reduce the computational effort required to perform the inversion \( z(\tau) \) implicit in the integral of (66) for very little loss in accuracy. It obviates the need for the large summations that appear in the first equality of (64) to be performed at every iteration of the complex-plane Newton-Raphson root finding.

The integration in (66) is equivalent to integration over the steepest descent path (SDP) of the phase function (69) and leads to a numerically exact result. On the other hand, depending on the accuracy required, the coefficients \( r_n \) of (65) may be used to construct the first few terms of the residuum series, of which the basic saddle-point (SP) approximation, given by (67), is the leading contribution.

One should note that, when this technique is used in a compound clutter problem, because the eigenvalues \( \gamma_m^{\infty} \) in (57) depend on the texture integration variable \( u \), a separate eigenvalue problem must be solved at each step of the numerical integration. We proceed to describe some approximation schemes that avoid this complication.

### B. Dalle Mese Giulì (DMG) Approximation

The generalized DMG approximation, as described in (12), is based on the use of simplified correlation matrices, chosen to reproduce the effective number of looks corresponding to the given correlation coefficients [2]. It has already been found to work well in the case of correlated gamma-fluctuating targets embedded in uncorrelated clutter speckle [12]. In that application, its adoption reduced computation time relative to the steepest-descent-path integration of the full effective model by a significant factor (primarily as all the \( M \)-fold summations over pulses are eliminated). It also provided the basis for proposing the MGF of the effective model. Furthermore, as discussed above, in addition to being an approximation to the effective model, it is also a bone fide approximation to the first-principles model.

In the generalized DMG approximation, the target and clutter correlation matrices \( C_t, C_c \) are constructed such that they commute, and are therefore simultaneously diagonalizable. Consequently, the speckle MGF of (56) holds with
\[ a_{qm}(u) = (1 - q + qu \gamma_{qm}^{\infty}) / M, \]
\[ a_{m}(u) = a_{qm}(u) + S \gamma_{qm}^{\infty} / (\kappa M), \tag{73} \]
where
\[ \{\gamma_{m}^{\infty} : m = 1, 2, \ldots , M\} = \text{eig}(C_t), \]
\[ \{\gamma_{m}^{s} : m = 1, 2, \ldots , M\} = \text{eig}(C_c). \tag{74} \]
Specifically,
\[ \gamma_{m}^{\infty} = 1 - \rho_{t} + M \rho_{t} \delta_{m} M, \]
\[ \gamma_{m}^{s} = 1 - \rho_{s} + M \rho_{s} \delta_{m} M, \tag{75} \]
Detection probability

where \( \delta_{m,n} \) denotes the Kronecker delta. The eigenvalues are expressed in terms of effective correlation coefficients, given by

\[
\rho'_s = \sqrt{(M/B - 1)/(M - 1)}, \\
\rho'_c = \sqrt{(M/L - 1)/(M - 1)},
\]

where the quantities \( B, L \) represent the effective numbers of looks on the target and clutter, respectively. These, in turn, can be obtained from the true correlation coefficients associated with an assumed correlation structure. For Kanter’s Gauss-Markov correlation model, one has

\[
B = M \cdot \left[ 1 + \frac{2\rho'_s^2}{1 - \rho'_s^2} \left( 1 - \frac{1}{M} \frac{1 - \rho'_c^2}{1 - \rho'_c^2} \right) \right]^{-1},
\]

\[
L = M \cdot \left[ 1 + \frac{2\rho'_c^2}{1 - \rho'_c^2} \left( 1 - \frac{1}{M} \frac{1 - \rho'_s^2}{1 - \rho'_s^2} \right) \right]^{-1}.
\]

All of this ensures that the DMG approximation reproduces the power variance of the underlying model to the extent possible within an effective-looks approach.

From (75), we see that \( c_m = \bar{c} \) and \( c_{qm} = \bar{c}_q \) for all \( m = 1, 2, \ldots, M - 1 \), so that the \( \tau \)-phase function reads

\[
\tau(z) = z + \ln(1 - c_0 z) \\
+ \kappa \left[ (M - 1) \ln(1 - \bar{c} z) + \ln(1 - c_M z) \right] \\
- (\kappa - 1) \left[ (M - 1) \ln(1 - \bar{c}_q z) + \ln(1 - c_{qM} z) \right].
\]

(77)

No Padé approximation is necessary in this case, as \( \tilde{r}_n = 0 \) for all \( n = 1, 2, \ldots \). The SP equation reads

\[
v = \frac{1}{s_0} + \kappa \left[ \frac{(M - 1)a_1(u)}{1 + a_1(u)s_0} + \frac{a_M(u)}{1 + a_M(u)s_0} \right] \\
- (\kappa - 1) \left[ \frac{(M - 1)a_{q1}(u)}{1 + a_{q1}(u)s_0} + \frac{a_{qM}(u)}{1 + a_{qM}(u)s_0} \right].
\]

(78)

Its solution is equivalent to finding the roots of a quintic polynomial. We may also note that, here,

\[
r_n = c^n_0 + \kappa \left[ (M - 1)c^n + c^n_{qM} \right] - (\kappa - 1) \left[ (M - 1)c^n_q + c^n_{qM} \right],
\]

for \( n = 1, 2, \ldots \), with the identity \( r_1 = 1 \) following from the SP equation.

Fig. 1 shows the detection probability for a false-alarm probability of \( P_{FA} = 10^{-n_{FA}} \), \( n_{FA} = 6 \) as a function of SIR for both exact effective model and DMG approximation, each computed via integration along the SDP and with the basic SP approximation. The difference between the SDP integration and SP approximation (solid versus dashed lines) is barely perceptible.

C. The Diagonal Approximation

Unlike the DMG approximation of the previous section, which has a theoretical basis arising from the effective looks concept of dealing with correlation, the diagonal approximation that we now introduce is an ad-hoc scheme, albeit one that turns out to be remarkably accurate, while conceptually simple, and exerting a somewhat reduced computational load relative to the full effective model. It comprises adoption of the effective model while pretending that the clutter and target correlation matrices commute, and can therefore be represented by their diagonalized forms. Then, as with the DMG approximation, we set

\[
a_{qm}(u) = (1 - q + qu\gamma^c_m)/M,
\]

\[
a_m(u) = a_{qm}(u) + S\gamma^s_m/(\kappa M),
\]

where

\[
\{\gamma^c_m : m = 1, 2, \ldots, M \} = \text{eig}(C_c),
\]

\[
\{\gamma^s_m : m = 1, 2, \ldots, M \} = \text{eig}(C_s),
\]

but now \( C_c, C_s \) are the actual correlation matrices rather than their simplified DMG counterparts.

The advantage of this over the full effective model is that, in the presence of compound clutter, eigenvalues of a target-plus-clutter correlation matrix do not have to be recomputed separately for each texture integration quadrature node. The diagonal approximation is manifestly exact for the following special cases:

- when \( \rho_c = 0 \) or \( \rho_a = 0 \);
- when \( \rho_c = \rho_a \), provided both the target and clutter have the same correlation structure;
- when \( M = 2 \) and in the \( M \to \infty \) limit;
- and obviously when \( S = 0 \) or \( q = 0 \).

It exhibits high accuracy across the full range of parameters. The deviation with respect to the exact result increases with power level \( v \), but lies well within 1\% for survival probabilities above \( 10^{-3} \). This is illustrated in Fig. 2 for a case where the deviation is amplified.

The magenta curve shows the difference in base-10 logarithms of the survival function between the exact effective model and the diagonal model. The green curve is the corresponding difference between the exact effective model and its basic saddle-point approximation. It can be seen that the basic SP approximation is worst around the mean of the distribution \( v = \bar{v} \), which occurs at a value of the power level normalized to mean interference given by \( \bar{v} = 1 + S = 6 \). The diagonal approximation slowly degrades with increasing power level \( v \),
but is generally superior. The power-axis interval shown in the graph spans a range of survival probabilities from unity to $10^{-3.5}$, as can be seen from the dashed blue curve representing the exact negative log-survival function.

D. Effective Model for Non-Fluctuating Targets

To derive the large-$\kappa$ limit of the effective model, we can look at

$$
\frac{\det(1/s + \sigma_n^2 + \sigma_C^2)}{\det(1/s + \sigma_n^2 + \sigma_C^2 + \sigma_C^2/\kappa)} \approx \frac{1}{\det(1 + \sigma_C^2 Q^{-1}(s)C_C/\kappa)} \sim \kappa^{-\sigma_C^2 Tr(C_C Q^{-1}(s))},
$$

from which we infer that

$$
\mathcal{M}_Z(s) \sim \kappa^{-\sigma_C^2 Tr(C_C Q^{-1}(s))}, \quad \kappa \to \infty.
$$

Equivalently, including the texture variable,

$$
\mathcal{M}_{Z_{sp}}^{\text{pk}}(s, u) \sim \exp \left[ -\text{Tr}\left\{ \ln(1 + (s/M)(1 - q + quC_C)) + (s/M)\left[ \frac{1}{\kappa} \sigma_C^2 Tr(C_C Q^{-1}(s)) \right] \right\} \right].
$$

We may now write $C_C = R_C^{-1} A_R R_C$, so that

$$
\text{Tr} \frac{C_C}{1 + (s/M)(1 - q + quC_C)} = \frac{M}{1 + (s/M)(1 - q + quC_C)},
$$

and introduce the quantities

$$
b_m = (1/M) \left[ R_C C_C R_C^{-1} \right]_{m,m} = (1/M) \sum_{n=1}^{M} \left[ R_C R_C^T \right]_{m,m} \gamma_n.
$$

This leads to the representation

$$
\mathcal{M}_{Z_{sp}}^{\text{pk}}(s, u) \sim \exp \left\{ -\sum_{m=1}^{M} \left[ \ln(1 + a_m(u)s) + S \cdot \frac{b_m s}{1 + a_m(u)s} \right] \right\},
$$

and it follows directly from (86) that $a_m = 1$ and that $b_m \geq 0$ for all $m = 1, 2, \ldots, M$. For a true steady target, in which case $C_C = \Omega$, where $\Omega_{ij} \equiv 1$ for all $i, j$, we obtain

$$
b_m = \frac{1}{M} \left( \sum_{n=1}^{M} [C_C]_{mn} \right)^2.
$$

A Swerling-0 target in the first-principles model (given the same target correlation matrix) yields, in (11), the expectation value

$$
\left\langle e^{-iu X_s} \right\rangle_{X_s} = \prod_{m=1}^{M} \cos(\sigma_s u_m),
$$

which can be shown to yield the MGF

$$
\mathcal{M}_Z(s) = \frac{1}{s^M \det Q(s)} e^{-\sigma_C^2 Tr(Q^{-1}(s))} \times \prod_{i,j=1, i\neq j}^{M} \cosh(\sigma_s^2 Q_{ij}^{-1}(s)).
$$

This result provides evidence that the effective and first-principles models are not strictly identical. However, as discussed in the Appendix, the fully phase-un correlated steady target is an exceptional case. In the general partially correlated case, the result generalizes to

$$
\mathcal{M}_Z(s) = \frac{1}{s^M \det Q(s)} e^{-\sigma_C^2 Tr(Q^{-1}(s))} \times \prod_{i,j=1, i\neq j}^{M} \cosh(\sigma_s^2 [L_i Q^{-1}(s) L_i^T]_{ij}),
$$

which follows from the representation $X_s = \sigma_s L_i^T B$, where the $B_m$ are iid Rademacher RVs.

VI. MONTE CARLO SIMULATION

While the first-principles model, as defined in (43), seems analytically intractable, it is highly amenable to Monte Carlo simulation, from which an empirical survival function can be derived. This enables a comparison with the effective model whose survival function can be explicitly computed via saddle-point techniques. Therefore, we can call upon the Kolmogorov-Smirnov statistic to test the null hypothesis that the empirical survival function derived from MC simulation
of the first-principles model and the analytically computed survival function for the effective model represent the same distribution.

A Monte Carlo (MC) simulation of the first-principles model, defined by (43), can be easily performed at the quadrature component level. In this case, the returned power RV is realized according to

\[
Z_{\text{avg}} / \sigma^2 = \frac{1}{2M} \sum_{m=1}^{M} \sum_{\eta=1}^{L_{\eta}} \left( \sqrt{1 - qH_{\alpha,\eta,m}} + \sqrt{qU\hat{X}_{c,\eta,m} + \sqrt{S\hat{X}_{s,\eta,m}}} \right)^2 .
\] (93)

Assuming Gauss-Markov correlated clutter, we can write

\[
\hat{X}_{c,\eta,m} = \rho_c \hat{X}_{c,\eta,m-1} + \sqrt{1 - \rho_c^2} H_{c,\eta,m} ,
\] (94)

for \( m = 2, 3, \ldots, M \), with \( \hat{X}_{c,\eta,1} = H_{c,\eta,1} \). Here, \( H_{\alpha,\eta,m} \sim \mathcal{N}(0,1) \) for both \( \alpha = n, c \). For the target RVs, we have

\[
\hat{X}_{s,\eta,m} = \sum_{m' = 1}^{M} \hat{Y}_{s,\eta,m'}[L_{\eta}][m/m'] ,
\] (95)

\[
\hat{Y}_{s,\eta,m} = (2B_{\eta,m} - 1) \sqrt{G_{\eta,m}} ,
\]

where the \( B_{\eta,m} \) represent iid Bernoulli trials such that \( B_{\eta,m} \sim \text{Bin}(1, 1/2) \), and the \( G_{\eta,m} \) are iid gamma variates such that \( G_{\eta,m} \sim \Gamma(\kappa/2, 2/\kappa) \). The RV \( U \) generates the clutter texture distribution. For K-distributed surface clutter, \( U \sim \Gamma(\nu, 1/\nu) \).

With the aid of MC simulation, we can compare the empirical survival function for the first-principles model with the computed survival function for the effective model. Fig. 3 compares two MC runs of the first-principles model for a gamma-fluctuating target with \( \kappa = 2 \) with the effective and DMG models each computed via integration along the SDP and basic SP approximation. A high level of coincidence is observed among all the cases. Fig. 4 draws the same survival function on a logarithmic scale where some discrepancies become apparent for the DMG approximation (red curve) in the tail. The dashed green and red curves for the effective model and DMG approximation arising from the SP approximation coincide with their SDP counterparts to within the line-width of the graph. Also, no difference between the first-principles (black and blue curves) and effective (green curve) models is evident. The MC sample size for each run was \( N_M = 10^4 \).

The null hypothesis that the two models generate the same distribution can be tested by means of the one-sample Kolmogorov-Smirnov (KS) statistic [18]. Since the population of returned powers associated with the first-principles model is sampled by MC simulations, we are able generate repeated samples, and thus construct an empirical distribution for the KS statistic. The results for the empirical survival function of the KS statistic are presented in Fig. 5 and Fig. 6 for target fluctuation classes \( \kappa = 1, 2 \), respectively.

We find that we are unable to reject the null hypothesis that the first-principles and effective models generate the same distribution at the \( \alpha = 0.01 \) significance level for almost any combination of parameters, with the exception of \( \kappa > 6 \) whenever \( C_n \) is very close (but not exactly equal) to the identity matrix, in which case a small discrepancy is detectable. The green region is generated by empirical survival functions derived via bootstrap resampling, and roughly delineates the 1% confidence interval. The red curve is the Dvoretzky-Kiefer-Wolfowitz (DKW) bound. The appearance of white space between the red curve and the lower edge of the green bootstrap region indicates that the null hypothesis should be rejected. The magenta curve is the survival function for the Kolmogorov distribution that represents the theoretical distribution for the KS statistic under the null hypothesis, and is generated by the stochastic process known as the Brownian bridge. We note that the red DKW bound essentially tracks the Brownian bridge due to the large sample size. We have also explicitly indicated the upper and lower confidence limits.

---

Fig. 3. Survival function for a \( \kappa = 2 \) correlated target in correlated K-clutter according to the various computational schemes. All curves except for those from the DMG approximation (red) coincide to within the line-width of the graph.

Fig. 4. Log-survival function for a \( \kappa = 2 \) correlated target in correlated K-clutter according to the various computational schemes. For the effective model (green) and DMG approximation (red), the dashed curve for the SP approximation coincides with the solid SDP curve to within the line-width of the graph.
Fig. 5. Empirical KS survival function for a $\kappa = 1$ correlated target in correlated K-clutter, comparing the first-principles and effective models.

Fig. 6. Empirical KS survival function for a $\kappa = 2$ correlated target in correlated K-clutter, comparing the first-principles and effective models.

Fig. 7. Empirical KS survival function for a $\kappa = 2$ correlated target in correlated K-clutter, comparing the first-principles model with the DMG approximation.

Statistical power generally increases with sample size, and is a measure of the sensitivity of the statistical test.

In order to quantify the power (i.e. sensitivity) of the KS test employed in the foregoing section, one may ask how much must one deform the distribution of the effective model away from its true functional form before the null hypothesis that it agrees with the empirical distribution of the first-principles model is rejected at the desired confidence level. In this section, we show that the required amount of deformation is tiny. In other words, if the survival function of the effective model were only very slightly different, then a discrepancy with respect to the first-principles model would be detected.

Thus, in order to characterize the statistical power of the KS test, we shall consider, as an alternative hypothesis, the slightly perturbed effective-model survival function [19], [20]

$$G(x) \equiv [\bar{F}(x)]^{1+\epsilon},$$

where $|\epsilon| \ll 1$. The value of the perturbation parameter $\epsilon$ can be related to the maximum difference $\delta_{\text{max}}$ between $\bar{F}(x)$ and $G(x)$. Writing

$$\delta(x) \equiv \bar{F}(x) - [\bar{F}(x)]^{1+\epsilon},$$

we obtain the derivative

$$\delta'(x) = f(x) \left[(1 + \epsilon)\bar{F}(x)^\epsilon - 1\right].$$

where $f(x)$ denotes the PDF associated with $\bar{F}(x)$. The extremum condition $\delta'(x) = 0$ is satisfied when

$$\bar{F}(x) = \left(\frac{1}{1 + \epsilon}\right)^{1/\epsilon}.$$

Hence, we find that

$$\delta_{\text{max}} = \bar{F}(x) [1 - \bar{F}(x)^\epsilon] = \epsilon \left(\frac{1}{1 + \epsilon}\right)^{1+1/\epsilon}.$$

This may be expressed as

$$\ln \delta_{\text{max}} = \ln \epsilon - (1 + 1/\epsilon) \ln(1 + \epsilon) = \ln \epsilon - 1 + O(\epsilon^2),$$

A. Statistical Power

False alarms occur in cases where the null hypothesis is rejected when it should not have been. Missed detections occur in cases where the null hypothesis is not rejected when it should have been. The power of a statistical test is formally defined as one minus the probability of a missed detection. The

---

As computed from Greenwood’s formula for a significance level of $\alpha = 0.01$, indicated by the dashed black curves. These curves are seen to bound the green bootstrap region.

We have also compared the first-principles model with the DMG approximation. The results for the empirical KS survival function for a $\kappa = 2$ target are presented in Fig. 5. Even though the DMG approximation agrees well with both the effective and first-principles model, as observed in the previous graphs, our statistical test has more than sufficient power to unambiguously reject the null hypothesis that it corresponds to the same distribution.

---

$\nu$ level of correlated $K$-clutter, comparing the first-principles and effective models.
The upshot of this is that while the KS test is able to detect the difference between the effective model and a deformation of it that is barely discernible, as shown in Fig. 8, it is unable to detect a discrepancy between the effective and first-principles models at the same significance level. It follows that the effective model is a sound proxy for the first-principles model.

VII. DISCUSSION

The first-principles and effective models always exhibit the same mean and variance, and are in exact agreement in the following situations:

- Uncorrelated clutter ($C_c = \mathbb{I}$);
- Gaussian target ($\kappa = 1$) with arbitrary target and clutter correlations;
- Swerling-0 target ($\kappa \to \infty$, $C_s = \Omega$) with arbitrary clutter correlation;
- Fully correlated target ($C_s = \Omega$) with arbitrary clutter correlation;
- $M \leq 2$ and $M \to \infty$;
- Commuting target and clutter correlation matrices (provided $C_s \neq \mathbb{I}$);
- Uncorrelated target ($C_s = \mathbb{I}$, (provided one sets $R_s = R_c$ when $\kappa > 1$);
- And, trivially, as $S \to 0$ and $S \to \infty$.

The first and final five cases apply to targets in all finite Swerling fluctuation classes $\kappa = 1, 2, \ldots$.

In all other cases studied, provided $\kappa \leq 6$ when $C_s$ is close to the identity matrix, the first-principles and effective models are seen to agree for all practical purposes, as demonstrated by examining the KS statistic. Thus, the effective model is a good proxy for the first-principles model, since the former lends itself to efficient computation via the saddle-point technique, whereas for the latter one must resort to MC simulation.

By studying the $\kappa \to \infty$ limit, we are able to show that the first-principles and effective models do not always coincide. The worst-case scenario occurs when $\rho_s = 1$, $\rho_c = 0^+$ and $q = 1$, with the SIR and integrated pulses around $S = 5$ and $M = 10$, respectively. Physically, this represents a highly unlikely scenario. The fluctuation parameter $\kappa$ must exceed $\kappa = 6$ before we are able to reject the null hypothesis.

In order to derive the first-principles MGF for the worst case discrepancy, we set $q = 1$, $C_s = \mathbb{I}$ and

$$C_c = \Omega = R_c^T \Lambda_c R_c, \quad \Lambda_c = \text{diag}(M, 0, \ldots, 0),$$

(102)

so that

$$M_{Z_{\alpha \nu}}(s) = \frac{1}{\det[sQ(s)]} e^{-(S/M) \text{Tr}(Q^{-1}(s))},$$

$$\times \prod_{i,j=1}^M \cosh\left(\frac{(S/M)Q_{ij}^{-1}(s)}{2}\right),$$

(103)

with $Q(s) = 1/s + \Omega/M$. Then, we have

$$Q^{-1}(s) = R_c^T \text{diag}(s/(1+s), s, \ldots, s) R_c$$

$$= s\mathbb{I} - \frac{s^2}{1+s} \Omega/M,$$

(104)
in which case
\[ \text{Tr}\{\tilde{Q}^{-1}(s)\} = (M - 1)s + s/(1 + s), \]
and so
\[ M_{Z_{\text{eff}}}(s) = \frac{1}{1 + s} e^{-(1-1/M)S_s} e^{-(S/M)s/(1 + s)}, \]
\[ \times \cosh^{M(M-1)} \left( \frac{(S/M^2)s^2}{1 + s} \right). \]

On comparing this with (83) we see that, by neglecting the \( \cosh \) term, the survival function for the corresponding effective model is recovered and it reduces to
\[ F_{\text{eff}}(v; S, M) = F_{\text{SW0}}(v - (1 - 1/M)S_s; S/M, 1), \]
where \( F_{\text{SW0}}(v; S, M) \) denotes the \( M \)-pulse Swerling-0 (i.e., Rician) survival function for SIR \( S \), normalized such that
\[ \int_0^\infty dv F_{\text{SW0}}(v; S, M) = 1 + S. \]

This serves to quantify the extent of the discrepancy. The mean KS statistic in this worst case is around 0.04 with the discrepancy highly localized in the vicinity of the knee of the survival function at \( v = (1 - 1/M)S \) as shown in Fig. 10. The survival function for the effective model was computed from (107) which holds strictly in the limit \( \rho_s \to 0 \). If we select a tiny but finite value, such as \( \rho_s = 10^{-4} \), the KS statistic drops to 0.008, for which the null hypothesis cannot be rejected.

In Table I, we compare the run-time and accuracy of each of the methods for calculating the effective model: These comprise the full effective model, the DMG approximation and the diagonal approximation, each computed within the basic saddle-point approximation (SP) and with integration along the path of steepest descent (SDP). The run-time is that required to compute a grid of 100 points on the survival function from the origin to the point where the probability has decreased to \( 10^{-3} \). Each probability error measures the maximum deviation of the survival function relative to the SDP result for the full effective model, expressed as both absolute and relative error. Each run-time and error represent an average of 100 randomized parameter combinations. The SIR is drawn in the range \( 1 \leq S \leq 10 \), the shape parameter is drawn in the range \( 1 \leq \nu \leq 10 \), and the clutter-to-interference ratio in the range \( 0.5 \leq q \leq 1 \). The remaining parameters are kept constant with values \( M = 100, \rho_c = 0.75, \rho_s = 0.95, \kappa = 2 \). The calculations were performed in MATLAB on a laptop computer with an Intel Core i5-4210U processor running at 1.70 GHz. We see that the DMG approximation is very fast but is less accurate than the diagonal approximation, which maintains a higher degree of accuracy but runs only marginally faster than the full effective model. The main draw-card of the diagonal model is its conceptual simplicity.

VIII. CONCLUSIONS

Our main conclusion is that, for all practical purposes, the effective model, defined in terms of its MGF, solves the problem of non-coherent detection of a gamma-fluctuating target in the presence of simultaneous arbitrary correlation of target RCS and clutter speckle, and solves it in a computationally efficient manner. One open question that arises from this work is whether there exists an alternative first-principles model whose MGF is exactly that of the effective model.

ACKNOWLEDGMENT

I would like to thank Stephen Bocquet and Luke Rosenberg for useful discussions and encouragement during the course of this research.

APPENDIX

A shortcoming of the first-principles model appears when dealing with an uncorrelated target, for which \( C_s = \mathbb{I} \). Here, \( \Lambda_s = \mathbb{I} \), and so \( L_s = R_s \) can be an arbitrary rotation matrix. However, the MGF given in Section IV is not independent of \( R_s \). A practical work-around is to select \( R_s = R_s \) in this case. This also serves to bring it into agreement with the associated effective model, introduced in Section V, which is explicitly independent of \( R_s \) for an uncorrelated target.

It may also appear that the first-principles model is not well-defined since the matrix \( L_s = \sqrt{\Lambda_s} R_s \) is not uniquely constructed from the eigenvalue decomposition \( C_s = R_s^T \Lambda_s R_s \). While this is true for \( C_s = \mathbb{I} \), it is not generally so: The uncorrelated target is an exceptional case. The underlying theory assumes a non-degenerate eigenvalue spectrum (which

---

**TABLE I**

| Method             | Run-time (sec) | Probability error |
|--------------------|----------------|------------------|
|                    | Absolute      | Relative         |
| Effective – SDP    | 5.59          | 3.58 x 10^{-2}   | 2.39 x 10^{-3} |
| Effective – SP     | 1.45          | 3.43 x 10^{-2}   | 2.30 x 10^{-3} |
| DMG – SDP          | 0.94          | 3.43 x 10^{-2}   | 2.44 x 10^{-3} |
| DMG – SP           | 0.66          | 6.75 x 10^{-2}   | 5.50 x 10^{-5} |
| Diagonal – SDP     | 5.13          | 1.92 x 10^{-4}   | 5.50 x 10^{-5} |
| Diagonal – SP      | 1.12          | 3.58 x 10^{-2}   | 2.44 x 10^{-3} |
is violated for an uncorrelated target). The eigenvalue decomposition partitions the Euclidean space $\mathbb{R}^M$ into $M$ orthogonal directions, and the columns of the matrix $R_c$ are the respective unit vectors associated with these directions. Each of these is uniquely determined up to an overall sign. Thus, we also have $C_\kappa = R_c^T \Lambda R_c$ where $R_c = \text{BR}_c$ for any diagonal matrix $B$ with diagonal elements $\pm 1$. Fortunately, the first-principles model as given in (43) is invariant under the transformation $L_1 \mapsto B L_1$.

It is worth pointing out that the first-principles MGF can be evaluated explicitly in the case of a fully correlated target ($C_\kappa = \Omega$, where $\Omega_{ij} = 1$ for all $i, j$) for any $\kappa$. Here, the matrix $L_1$ is such that, for any matrix $A$,

$$L_1 A^{T} = \text{diag} \left( \text{Tr} \{ \Omega A \}, 0, \ldots, 0 \right).$$  \hspace{1cm}  \text{(109)}$$

In particular,

$$L_1 C_\kappa^{-1} s^{T} L_1 = \text{diag} \left( \text{Tr} \{ C_\kappa^{-1} (s) \}, 0, \ldots, 0 \right),$$  \hspace{1cm}  \text{(110)}$$

in which case

$$\mathbf{Y}^{T} (L_1 C_\kappa^{-1} s^{T} L_1) \mathbf{Y} = \text{Tr} \{ C_\kappa^{-1} (s) \} \cdot \hat{Y}_1^2.$$  \hspace{1cm}  \text{(111)}$$

It follows that

$$M_Z(s) = \frac{1}{s^M |\text{det} Q(s)|} \left( \text{Tr} \{ C_\kappa^{-1} (s) \} \cdot \hat{Y}_1^2 \right)^2.$$  \hspace{1cm}  \text{(112)}$$

The expectation is given by

$$\left\langle \exp \left( -\frac{s^2}{2} \text{Tr} \{ C_\kappa^{-1} (s) \} \cdot \hat{Y}_1^2 \right) \right\rangle_{\hat{Y}_1} = \frac{1}{\Gamma(\kappa/2)} \int_0^{\infty} dx x^{\kappa/2 - 1} \times \exp \left\{ - \left[ 1 + (\sigma^2_0/\kappa) \text{Tr} \{ C_\kappa^{-1} (s) \} \right] x \right\}$$

$$= \left[ 1 + (\sigma^2_0/\kappa) \text{Tr} \{ C_\kappa^{-1} (s) \} \right]^{-\kappa/2}.$$  \hspace{1cm}  \text{(113)}$$

Next, let us look at

$$\ln \text{det} \left[ I + (\sigma^2_0/\kappa) C_\kappa^{-1} (s) \right] = \text{Tr} \left\{ \ln \left[ I + (\sigma^2_0/\kappa) C_\kappa^{-1} (s) \right] \right\}$$

$$= \sum_{n=1}^{\infty} \frac{1}{n} \left( \sigma^2_0/\kappa \right)^n \text{Tr} \{ C_\kappa^{-1} (s) \}^n,$$  \hspace{1cm}  \text{(114)}$$

and observe that

$$\text{Tr} \{ C_\kappa^{-1} \}^n = \text{Tr} \{ L_1 C_\kappa^{-1} L_1^T \}^n = \left( \text{Tr} \{ C_\kappa^{-1} \} \right)^n,$$  \hspace{1cm}  \text{(115)}$$

from which it follows that

$$M_Z(s) = \frac{1}{s^M |\text{det} Q(s)|} \left( \text{det} \left[ I + \sigma^2_0 C_\kappa^{-1} (s)/\kappa \right] \right)^{-\kappa},$$  \hspace{1cm}  \text{(116)}$$

which, as one can see by appealing to [49] or [83], is in agreement with the effective model discussed in Section V.
Josef Zuk was awarded a B.Sc.(Hons) in Physics and Mathematics from the University of Melbourne in 1982. In 1985, he completed a D.Phil. in Theoretical Particle Physics at the University of Oxford, where he remained as a research fellow until 1987, before moving on to the Max Planck Institute for Nuclear Physics in Heidelberg, Germany to take up a research fellowship in Theoretical Condensed-Matter Physics. In 1990, he commenced a faculty position in the Department of Physics at the University of Manitoba, Winnipeg, Canada. Having joined DST Group (then DSTO) in 1995, he is currently a senior research scientist in the Joint and Operations Analysis Division (JOAD), where he works in operations research, and is mainly engaged in operational aspects of radar performance modelling and simulation, particularly for airborne maritime surveillance radar systems.