

The Application of Time Series Modelling and Monte Carlo Simulation: Forecasting Volatile Inventory Requirements

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Abstract

During the assembly of internal combustion engines, the specific size of crankshaft shell bearing is not known until the crankshaft is fitted to the engine block. Though the build requirements for the engine are consistent, the consumption profile of the different size shell bearings can follow a highly volatile trajectory due to minor variation in the dimensions of the crankshaft and engine block. The paper assesses the suitability of time series models including ARIMA and exponential smoothing as an appropriate method to forecast future requirements. Additionally, a Monte Carlo method is applied through building a VBA simulation tool in Microsoft Excel and comparing the output to the time series forecasts.

Keywords

Forecasting, Time Series Analysis, Monte Carlo Simulation

1. Introduction

Inventory control is an essential element within the discipline of operations management and serves to ensure sufficient parts and raw materials are available for immediate production needs while minimising the overall stock holding at the point of production and throughout the supply chain. Methodologies including Materials Requirements Planning and Just-in-Time Manufacturing have evolved to manage the complexities of supply management supported by an extensive academic literature. Similarly extensive study into the inventory profiles for spare part and service demand has also been prevalent over the last 50 years.

This paper presents a study of an inventory process that does not fit into the standard inventory models within conventional operations management or for service and spare parts management. In high volume internal com-

bustion engine manufacturing, the demand profile for crankshaft shell bearings follow a highly variable demand profile though there is consistent demand for the engines. The paper reviews the application of both time series analysis and a Monte Carlo simulation method to construct a robust forecast for the shell usage consumption. Section 2 presents the problem statement. Time series analysis is reviewed in Section 3. The Monte Carlo simulation method written in Microsoft Excel VBA is presented in Section 4. Forecasts generated by both the time series models and the simulation are assessed in Section 5 and concluding remarks are presented in Section 6. The analysis utilises usage data provided by a volume engine manufacturer over a 109 week production build period.

2. Problem Statement

The construction of petrol and diesel engines involve fitting a crankshaft into an engine block and attaching piston connecting rods to the crankshaft crank pins [1]. The pistons via the connecting rods turn the crankshaft during the power stroke of the engine cycle to impart motion. Shell bearings are fitted between the crank journals and crank pins and the engine block and connecting rods. During engine operation a thin film of oil under pressure separates the bearing surface from the crankshaft journals and pins. For the smooth operation of the engine, the thickness of the oil film has to be consistent for each crank journal and pin. Though the crankshaft, connecting rods and engine block are machined to exceptionally tight tolerances minor deviations in the dimensions of the machined surfaces mean that the thickness of the oil film will not be consistent across the individual crank journals and pins.

To overcome the problem, the engine designer specifies a range of tolerance bands within an overall tolerance for the nominal thickness of the shell bearing. Similar tolerance bands are specified for the machined surfaces of the engine block and crankshaft. A shell bearing whose thickness is defined within a specific tolerance band is identified by a small colour mark applied during the shell manufacturing process. The engine designer creates a “Fitting Matrix”, where the combination of the tolerance bands for the engine block and crankshaft are compared against which the appropriate shell bearing can be selected. During the engine assembly, the selection process is automated. Embossed marks on the crankshaft and engine block specify the tolerance bands of each machined surface. The marks are either optically scanned or fed into a device that creates a visual display of the colours of the shells to select for assembly. Selection is rapid and does not impede the speed of engine assembly. **Table 1** illustrates the set of tolerance bands and associated “Fitting Matrix” for the selection of main bearing shells for a typical engine.

Over time the usage profile for some shells thicknesses can show considerable variation against a consistent demand for the engine itself. The high variation poses a difficult procurement problem for engine assemblers that support high volume automotive production. It is difficult to determine what the usage consumption for high usage shells will be over a future time horizon as the shell bearings can be globally sourced and can require lead times of between 3 and 6 months.

Figure 1 illustrates the consumption trajectory of journal shell sizes over a 109-week sample period that has high demand profiles. The shells are identified by the colour green and yellow. The green shell consumption is highly volatile with sporadic spikes in demand. The yellow shell consumption after showing a rapid decline over

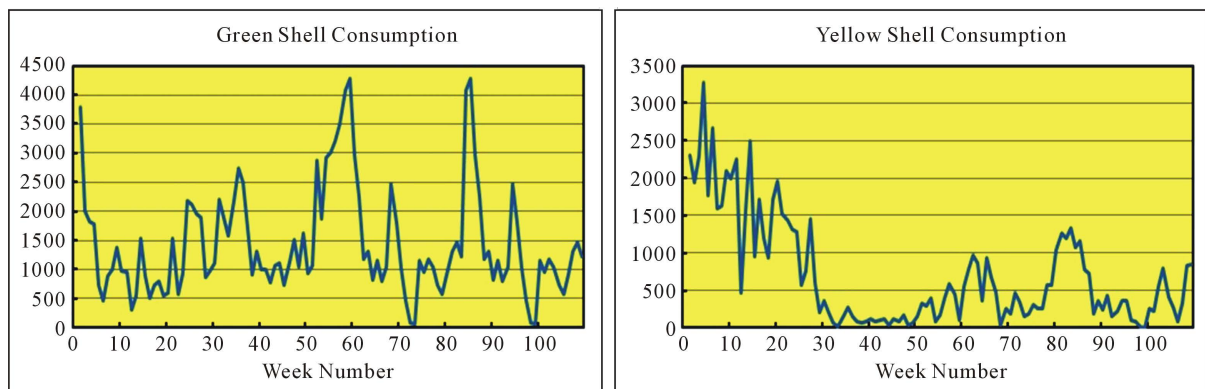


Figure 1. Time series trajectories for green and yellow shell consumption.

Table 1. Bearing selection-fitting matrix.

Dimension Table								
Engine Block: Main Bearing			Crankshaft Main Bearing			Main Bearing		
Dia = 59 mm	Upp Tol	+0.024 μm	Dia = 55 mm	Upp Tol	+0.008 μm	Thickness = 2 mm	Upp Tol	+0.012 μm
	Lwr Tol	-0 μm		Lwr Tol	-0.016 μm		Lwr Tol	-0.06 μm
Mark	Range (μm)		Mark	Range (μm)		Colour	Range (μm)	
A	0	+6	1	+8	+4	Blue	+12	+9
B	+6	+12	2	+4	0	Black	+9	+6
C	+12	+18	3	0	-4	Brown	+6	+3
D	+18	+24	4	-4	-8	Green	+3	0
			5	-8	-12	Yellow	0	-3
			6	-12	-16	Pink	-3	-6
Fitting Matrix								
Main Bearing								
		A	B	C	D			
Engine Block	1	Pink	Pink/Yellow	Yellow	Green			
	2	Pink/Yellow	Yellow	Green	Green/Brown			
	3	Yellow	Green	Green/Brown	Brown			
	4	Green	Green/Brown	Brown	Black			
	5	Green/Brown	Brown	Black	Black/Blue			
	6	Brown	Black	Black/Blue	Blue			

the first 30 weeks of the sample period settles to lower and less volatile demand than the green shell.

3. Time Series Analysis

Succinctly, a time series is a record of the observed values of a process or phenomena taken sequentially over time. The observed values are random in nature rather than deterministic where the random behaviour is more suitable to model through the laws of probability. Systems or processes that are non-deterministic in nature are referred to as *stochastic* where the observed values are modelled as a sequence of random variables. Formally:

A stochastic process $\{Y(t); t \in T\}$ is a collection of random variables where T is an index set for which all the random variables $Y(t)$, $t \in T$, are defined on the same sample space. When the index set T represents time, the stochastic process is referred to as a *time series*, [2].

While time series analysis has extensive application, the purpose of the analysis is two-fold [3]. Firstly to understand or model the random (stochastic) mechanism that gives rise to an observed series and secondly to predict or forecast the future values of a series based on the history of that series and, possibly, other related series or factors.

Time series are generally classified as either stationary or non stationary. Simplistically, stationary time series process exhibit invariant properties over time with respect to the mean and variance of the series. Conversely, for non stationary time series the mean, variance or both will change over the trajectory of the time series. Stationary time series have the advantage of representation by analytical models against which forecasts can be produced. Non stationary models through a process of differencing can be reduced to a stationary time series and are so open to analysis applied to stationary processes [4].

An additional method of time series analysis is provided by smoothing the series. Smoothing methods estimate the underlying process signal through creating a smoothed value that is the average of the current plus past observations.

3.1. Analysis of Stationary and Non-Stationary Time Series

Stationary time series are classified as having time invariant properties over the trajectory of the time series. It is these time invariant properties that are stationary while the time series itself appears to fluctuating in a random manner. Formally, an observed time series $\{Y(t); t \in T\}$ is *weakly stationary* if the following properties hold:

- 1) $E[Y(t)] = \mu, \forall t \in T$
- 2) $\text{Var}[Y(t)] = \sigma^2 < \infty$
- 3) $\text{Cov}(Y_t, Y_{t-j}) = \gamma_j, \forall t, j \in T$

Further, the time series is defined as strictly stationary if in addition to Properties 1-3, if subsequently:

- 4) The joint distribution of $\{Y(t_1), Y(t_2), \dots, Y(t_k)\}$ is identical to $\{Y(t_1+h), Y(t_2+h), \dots, Y(t_k+h)\}$ for any $t_i, h \in T$

Conversely, a non-stationary time series will fail to meet either or both the conditions of Properties 1 and 2.

Property 3 states that the covariance of lagged values of the time series is dependent on the lag and not time. Subsequently the autocorrelation coefficient ρ_k at lag k is also time invariant and is given by

$$\rho_k = \frac{\text{Cov}(y_t, y_{t+k})}{\text{Var}(y_t)} = \frac{\gamma_k}{\gamma_0} \quad (1)$$

The set of all $\rho_k, k = 0, 1, 2, \dots, n$ forms the Autocorrelation Function or ACF. The ACF is presented as a graphical plot. **Figure 2** provides an example of a stationary time series with the associated ACF diagram. Successive observations of a stationary time series should show little or no correlation and is reflected in the ACF plot showing a rapid decline to zero.

Conversely, the ACF of a non stationary process will show a gentler decline to zero. **Figure 3** replicates the ACF for a non-stationary random walk process.

Stationary time series models are well defined throughout the time series literature where a full treatment of their structure can be found. Representative references include [5]-[8].

A consistent view of a time series is that of a process consisting of two components a signal and noise. The signal is effectively a pattern generated by the dynamics of the underlying process generating the time series. The noise is anything that perturbs the signal away from its true trajectory [7]. If the noise results in a time series that consists of uncorrelated observations with a constant variance, the time series is defined as a white noise process. Stationary time models are always white noise processes where the noise is represented by a sequence of error or shock terms $\{\varepsilon_t\}$ where $\varepsilon_t \sim N(0, \sigma)$. The time series models applicable to stationary and non stationary time series are listed in **Table 2**.

From **Table 2**, it is seen that by setting the $p, d,$ and q parameters to zero as appropriate, the $\text{MA}(q), \text{AR}(p)$ and $\text{ARMA}(p, q)$ processes can be presented as sub processes of the $\text{ARIMA}(p, d, q)$ process and is illustrated in the hierarchy presented in **Figure 4**. The stationary processes can be thought of as ARIMA processes that do not require differencing.

Table 2. Description of time series models.

1) **Moving Average Process $\text{MA}(q)$:** The time series y_t is defined as the sum of the process mean and the current shock value plus a weighted sum of the previous “ q ” past shock values.

2) **Autoregressive Process $\text{AR}(p)$:** The time series y_t is presented as a linear dependence of weighted “ p ” observed past values summed with the current shock value and a constant.

3) **Autoregressive Moving Average $\text{ARMA}(p, q)$:** The time series y_t is presented as a mixture of both moving average and autoregressive terms. $\text{ARMA}(p, q)$ processes require fewer parameters when compared to the AR or MA process (Chatfield, 2006).

4) **Autoregressive Integrated Moving Average $\text{ARIMA}(p, d, q)$:** A non stationary time series is transformed into a stationary time series through a process of differencing. The ARIMA process differences a time series at most d times to obtain a stationary $\text{ARMA}(p, q)$ process.

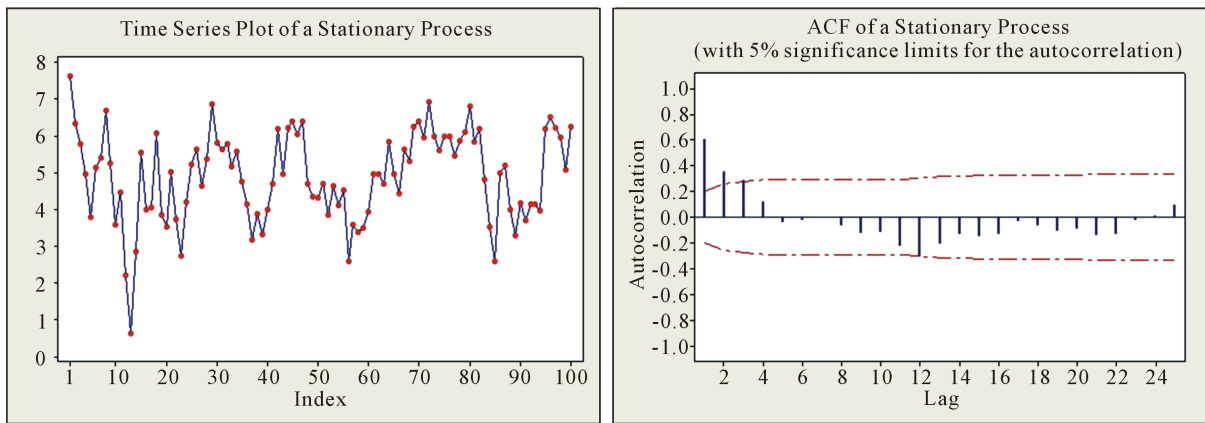


Figure 2. Stationary time series—example ACF.

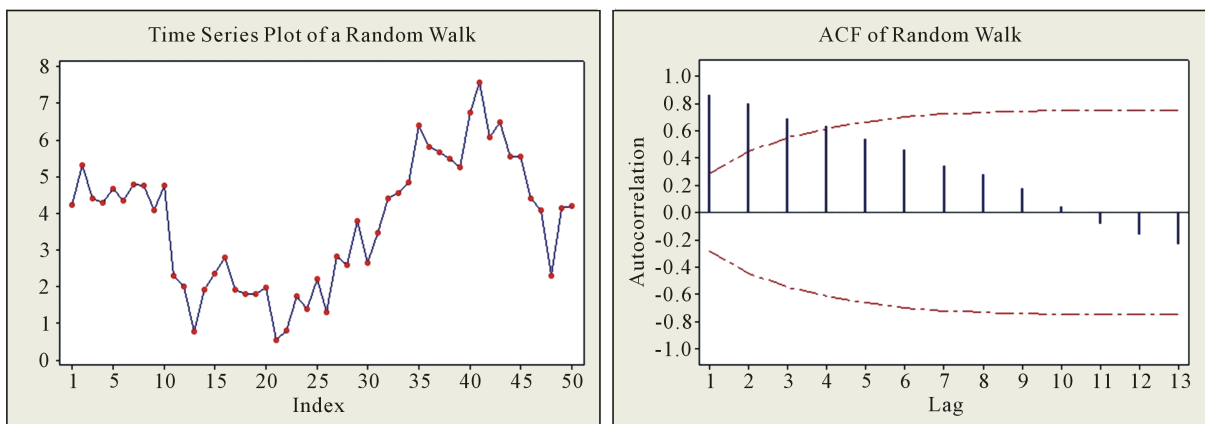


Figure 3. Non-stationary time series example ACF (Random walk).

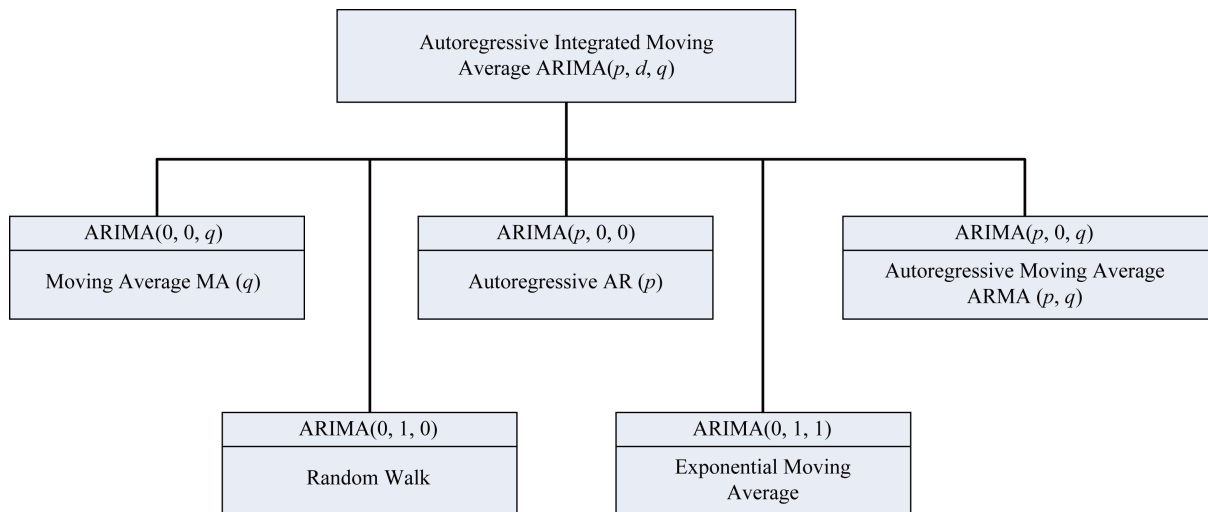


Figure 4. Time series model hierarchy.

Moreover, the ARIMA process reveals time series models that are not open to analysis by the stationary representations. Two such models are included in **Figure 4**, the random walk model, ARIMA(0, 1, 0) and the exponential moving average model, ARIMA(0, 1, 1).

Table 3 returns the modelling processes that are applied to stationary and non-stationary time series. The

Table 3. Synthesis of time series models.

Process	Model	Stationary Condition
MA(q)	$y_t = \mu + \varepsilon_t - \theta_1 \varepsilon_{t-1} - \theta_2 \varepsilon_{t-2} - \dots - \theta_q \varepsilon_{t-q}$ $y_t = \mu + \Theta(B) \varepsilon_t$	None: MA(q) process is always stationary.
AR(p)	$y_t = \delta + \phi_1 y_{t-1} + \phi_2 y_{t-2} + \dots + \phi_p y_{t-p} + \varepsilon_t$ $\Phi(B) y_t = \delta + \varepsilon_t$	$\sum_{i=1}^p \phi_i < 1$
ARMA(p, q)	$y_t = \delta + \sum_{i=1}^p \phi_i y_{t-i} + \varepsilon_t - \sum_{i=1}^q \theta_i \varepsilon_{t-i}$ $\Phi(B) y_t = \delta + \Theta(B) \varepsilon_t$	AR(p) component is stationary.
ARIMA(p, d, q)	$\Phi(B)(1-B)^d y_t = \delta + \Theta(B) \varepsilon_t$	The AR(p) component is stationary after the series is differenced d times.

equations of the models are replicated in their standard form and in terms of the backshift operator B (see Appendix 1). Time series models expressed in terms of the backshift operator are more compact and enable a representation for the ARIMA process that has no standard form equivalent.

3.2. Model Identification

Initially, inspection of the ACF of a time series is necessary to determine if the series is stationary or will require differencing. There is no analytical method to determine the order of differencing though empirical evidence suggests that generally first order differencing ($d = 1$) is sufficient and occasionally second order differencing ($d = 2$) should be enough to achieve a stationary series [7].

The ACF is also useful to indicate the order of the MA(q) process as it can be shown that for $k > q$, $\rho_k = 0$. Hence the ACF will cut off at lag q for the MA(q) process. The ACF of the AR(p) and ARMA(p, q) processes both exhibit exponential decay and damped sinusoidal patterns. As the form of the ACF's for these processes is indistinguishable, identification of the process is provided by the Partial Autocorrelation Coefficient (PACF).

The properties of the PACF are discussed extensively in the time series literature and in particular a comprehensive review provided in [9]. Descriptively, the PACF quantifies the correlation between two variables that is not explained by their mutual correlations to other variables of interest. In an autoregressive process, each term is dependent on a linear combination of the preceding terms. In evaluating the autocorrelation coefficient ρ_k , the term y_t is correlated to y_{t-k} . However, y_t is dependent on y_{t-1} which in turn is dependent on y_{t-2} and the dependency propagates throughout the time series to y_{t-k} . Consequently, the correlation at the initial lag of the time series propagates throughout the series. The PACF evaluates the correlation between x_t is and x_{t-k} through removing the propagation effect.

The PACF can be calculated for any stationary time series. For an AR(p) process, it can be shown that the PACF cuts off at lag p . The PACF for both the MA(q) and ARMA(p, q) process the PACF is a combination of damped sinusoidal and exponential decay.

The structure of a stationary time series is determined through inspection of both the ACF and PACF diagrams of the series. **Table 4** presents the classification of the stationary time behaviour (adapted from [7] [8]).

Neither the ACF nor PACF yield any useful information with respect to identifying the order of the ARMA (p, q) process. Though there are additional methods that can aid the identification of the required order of the process [7] [8] including the extended sample autocorrelation function (ESACF), the generalised sample partial autocorrelation function (GPACF), and inverse autocorrelation function (IACF) as suitable methods for determining the order of the ARMA model. However, with respect to investigating the structure of a time series both sets of authors agree that with the availability of statistical software packages it is easier to investigate a range of models with various orders to identify the appropriate model and forego the need to apply these additional methods.

The estimation of the parameter coefficients ϕ_i and θ_i can be estimated through Maximum Likelihood Estimation (MLE). An extensive explanation of the MLE application to estimate the parameters of each the models

Table 4. Classification of stationary time series behavior.

Model	ACF	PACF
MA(q)	Cuts off after lag q	Infinite exponential decay and/or damped sinusoid—tailing off
AR(p)	Infinite exponential decay and/or damped sinusoid—tailing off	Finite-cuts off after lag p
ARMA(p, q)	Infinite exponential decay and/or damped sinusoid—tailing off	Infinite exponential decay and/or damped sinusoid—tailing off

listed in **Table 2** is provided in [10]. Additionally the parameters of the AR(p) process can be calculated with the Yule Walker Equations [4] [9].

The robustness of a derived model is assessed through residual analysis of the model. For the ARMA(p, q) process the residual are obtained from

$$\hat{\varepsilon}_t = y_t - \left(\hat{\delta} + \sum_{i=1}^p \hat{\phi}_i y_{t-i} - \sum_{i=1}^q \hat{\theta}_i \varepsilon_{t-i} \right) \quad (2)$$

If the residual values exhibit white noise behaviour, the ACF and PACF diagrams of the residual values should not show any discernable pattern then the appropriate model has been chosen and the correct orders of p and q have been correctly identified.

3.3. Generation of Forecasts

Establishing a model that describes the structure of an observed time series enables meaningful forecasts to be drawn from the model. Forecasting methods for each of the models in **Table 2** are succinctly described in [11] and are defined as follows:

AR(p) Process:

The forecast for the AR(p) process is based on the property that the expectation of the error terms ε_i are zero. The forecast is produced iteratively from the last observation where **at each iteration a one step** ahead forecast is produced. The step ahead is denoted by

$$\begin{aligned} \tau = 1: \hat{y}_t(1) &= \delta + \phi_1 y_{t-1} + \phi_2 y_{t-2} + \dots + \phi_p y_{t-p+1} \\ \tau = 2: \hat{y}_t(2) &= \delta + \phi_1 \hat{y}_t(1) + \phi_2 y_{t-2} + \dots + \phi_p y_{t-p+2} \end{aligned}$$

At each successive iteration, the most recent observation drops out of the forecast and replaced by the previous forecast value. At $\tau > p$, each term is a forecasted value and continuing the iteration process, the forecast converges onto the mean of the AR(p) process.

MA(q) Process:

The expectation of the ε_i terms in the MA(q) process follow a white noise process and so the expectation of all future values of $\varepsilon_i, i > 0$ is zero. Hence for $t > q$ the forecast of a MA(q) process is just the mean value of the process, μ .

ARMA(p, q) Process:

The forecast for the ARMA(p, q) process is the combination of the results from the pure AR(p) and MA(q) processes. For $t > q$, the error terms completely drop out of the forecast.

3.4. Smoothing Methods

Smoothing methods provide a class of time series models that have the effect of reducing the random variation of a time series with the effect of exposing the process signal within the time series. A variety of smoothing models that average the series continually over a moving span of observed values or fit a polynomial approximation to a restricted interval of the series are presented in [12].

Of the smoothing techniques available, the exponential smoothing model has proved useful due to its simplicity of application. First order exponential smoothing is defined by the following recurrence relation:

$$\tilde{y}_t = \lambda y_t + (1 - \lambda) \tilde{y}_{t-1} \quad (3)$$

where λ is the smoothing factor ($0 < \lambda < 1$).

Effectively, first order exponential smoothing is a linear combination of the current observation plus the discounted sum of all previous observations due to the smoothing factor λ . Moving back through the recursive relation $(1 - \lambda)$ geometrically decays and so the older observations have a diminishing contribution to the smoothed estimation of the current value.

Though the recursive relation defined in Equation (3) can be expressed as an ARIMA(0, 1, 1) model, the method was initially developed from first principles in [13] as a means of forecasting inventory.

First order exponential smoothing will for trending data lag behind an increasing trend and lead a decreasing trend. Second order exponential smoothing overcomes this problem as does increasing the value of the smoothing factor [7]. However, providing the time series is showing no systematic trend, first order exponential smoothing is an adequate model to analyse a time series [6].

The forecast generated from first order exponential smoothing is just the value of the smoothed current value and in principle would extend over all future values. However, as more observed values become available it makes sense to update the forecast.

4. Simulation Application to Evaluate Consumption Rate (Monte Carlo Method)

Frequently in real world scenarios due to the complexity of the system under investigation it may not be possible to evaluate the systems behaviour by applying analytical methods. Under such conditions an alternative approach to model such system is through creating a simulation. Succinctly, simulation methods provide an alternative approach to studying system behaviour through creating an artificial replication or imitation of the real world system [14]. The applications where simulation methods may be useful is extensive and include diverse disciplines such as manufacturing systems, flight simulation, construction, healthcare, military and economics [15] [16].

Systems or processes that can be modelled through an underlying probability distribution are open to simulation through the Monte Carlo method [17]. The method simulates the behaviour of a system by taking repeated sets of random numbers from the underlying probability distribution of the process under investigation. The development of the method is attributed to the work of Stanislaw Ulam and John von Neumann during the late 1940's to aid their work in atomic physics for the development of nuclear weapons. Analytical methods were proving difficult if not impossible to apply and Ulam turned to random experimentation to elicit system attributes and behaviour [18].

A specific application of the Monte Carlo method is dependent on the nature of the underlying probability distribution of the system or process under investigation. However the method application is consistent and will follow the steps outlined in **Table 5**.

Application of Monte Carlo Method to Shell Consumption Case Study

Visual inspection of the time series trajectories of the shell consumption does not yield an obvious probability distribution so Stage 1 of the Monte Carlo process begins by examining the histograms of the observed series.

Table 5. Monte Carlo method process stages (Adapted from [17]).

Stage 1 Define a distribution of possible inputs for each input random variable: Requires recognition of the underlying probability distribution of the process. This may be directly apparent or may require empirical observation of the process under investigation.

Stage 2 Generate outputs randomly from those distributions: Requires the selection of an appropriate random number generator to model the observed probability distribution. Random number generators are generally available in most statistical software packages and Micro Soft Excel.

Stage 3 Perform a deterministic computation using that set of outputs: Involves computing the desired output variable or variables from the generated random numbers.

Stage 4 Aggregate the results of the individual computations into the final result: The aggregation process is dependent on the specific simulation but could be as straightforward as computing the average of the simulated results.

Figure 5 returns the histograms of the consumption of the Green and Yellow Shell consumption for the full trajectory of observed values.

The histograms for each shell are returned in **Figure 4** and indicate the data is negatively skewed and so drawing random numbers from a skewed normal distribution for the simulation is deemed appropriate. It can also be shown that the distributions for shorter time spans of the observed values are also skewed.

Stages 2 to 4 of the Monte Carlo process are embedded in a Visual Basic for Applications (VBA) program created in Microsoft Excel to carry out the simulation. **Figure 6** provides a schematic of the VBA simulation programme.

The VBA programme incorporates a skewed normal random number generation algorithm developed in [19] that generates random numbers based on the skew value of sample input data. The Excel spreadsheet contains the consumption history for each shell. Upon invoking the simulation, the required sample and forecast periods

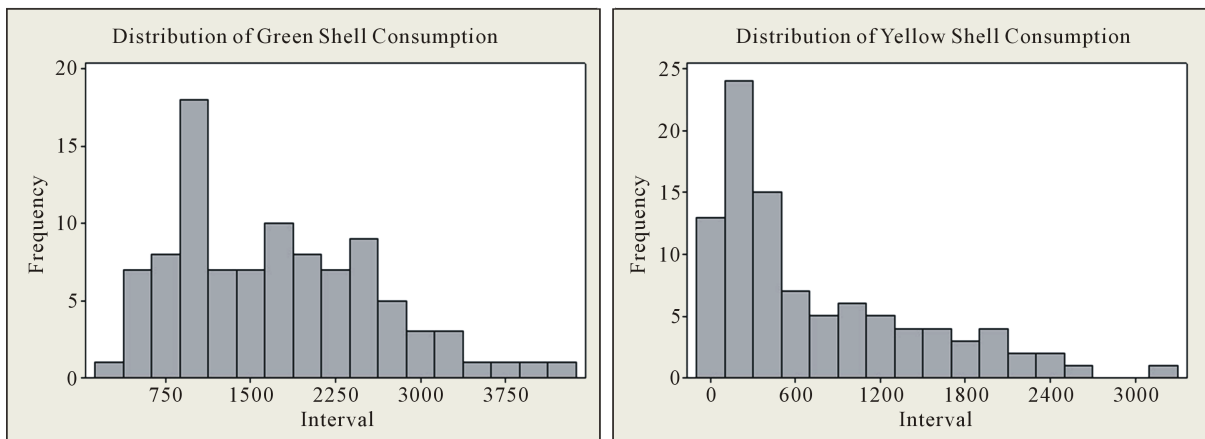


Figure 5. Histograms of green and yellow shell consumption.

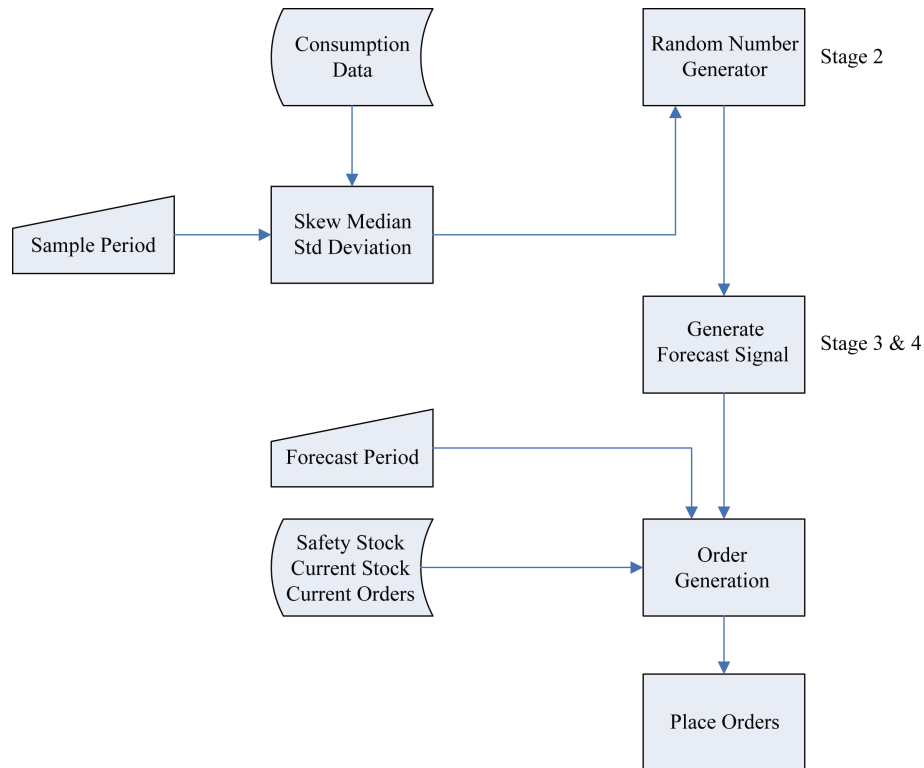


Figure 6. Schematic of VBA simulation programme.

are read from input data in the spreadsheet. For each shell, against the sample period, the skew, median and standard deviation values are calculated and input into the random number generator. For each week of the sample period, the generator calculates 100 positive random numbers to create an average simulated consumption for each week. Since it is impossible to have negative consumption, negative random numbers are rejected. Against each weekly average, an overall average is computed to create the forecast signal.

The random number generator satisfies Stage 2 of Monte Carlo process, while Stages 3 and 4 are satisfied through computation of the forecast signal.

Additional functionality is provided in the simulation tool that over the forecast period will generate orders taking into account current stock levels, safety levels and orders already placed.

5. Generation of Forecasts for the Bearing Shell Consumption

The time series and the Monte Carlo methods described in the preceding sections are applied to the historical shell bearing consumption usage to generate forecasts to create orders to satisfy future engine build requirements. The forecasts of each model are compared to determine if there is either a consistent or significant differences between each method.

5.1. ARIMA (p, d, q)

The purpose of the ARIMA(p, d, q) process is to establish the underlying model that describes the time series through specifying the p and q parameters once the appropriate order of differencing d has identified a stationary process. Model identification is primarily based on inspection of the ACF and PACF diagrams and referencing the Classification Table (Table 4). The robustness of the forecast generated from the identified model is determined by Residual Analysis where primarily if the ACF and PACF diagrams of the residual values show no discernible pattern, the forecast can be assessed as robust [2]-[11]. It is considered that the identification of a time series model requires both judgement and experience and it is recommended that the iterative model building process illustrated in Figure 7 is followed [8] [9].

The forecasting process is illustrated against the Green Shell consumption replicated in Figure 1. The analysis is carried out using the Minitab[®] statistical package. The ACF and PACF diagrams are replicated in Figure 8 and with reference to Table 4 indicates that the time series follows an AR(1) process.

Table 6 returns the output generated against 40 observations. The ACF and PACF diagrams of the residual values are returned in Figure 9 and do not show any discernible pattern. The Chi Square statistics appear to be

Table 6. Model estimation for green usage.

ARIMA Model: Green				
Final Estimates of Parameters				
Type	Coef	SE Coef	T	P
AR 1	0.6940	0.1168	5.94	0.000
Constant	368.1	104.8	3.51	0.001
Mean	1203.0	342.5		
Number of observations: 40				
Residuals:	SS = 16,691,055 (backforecasts excluded)			
	MS = 439,238; DF = 38			
Modified Box-Pierce (Ljung-Box) Chi-Square statistic				
Lag	12	24	36	8
Chi-Square	14.9	21.1	23.9	*
DF	10	22	34	*
P-Value	0.135	0.518	0.902	*

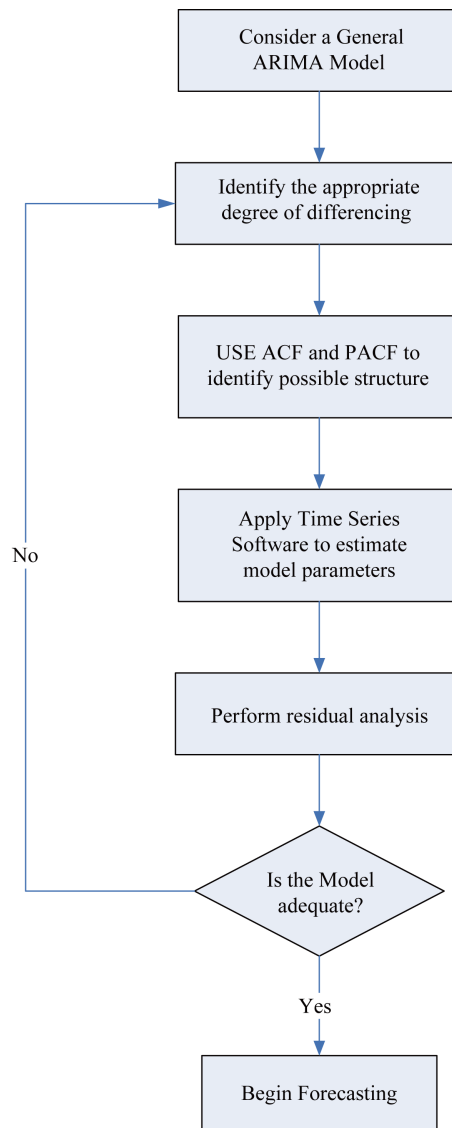


Figure 7. ARIMA model building stages (Adapted from [8] [9]).

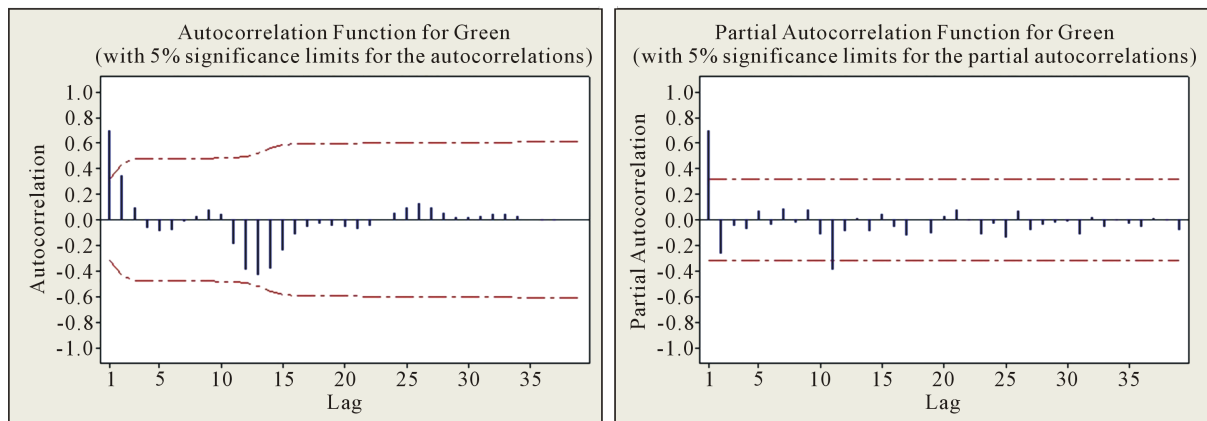


Figure 8. ACF and PACF of green shell usage.

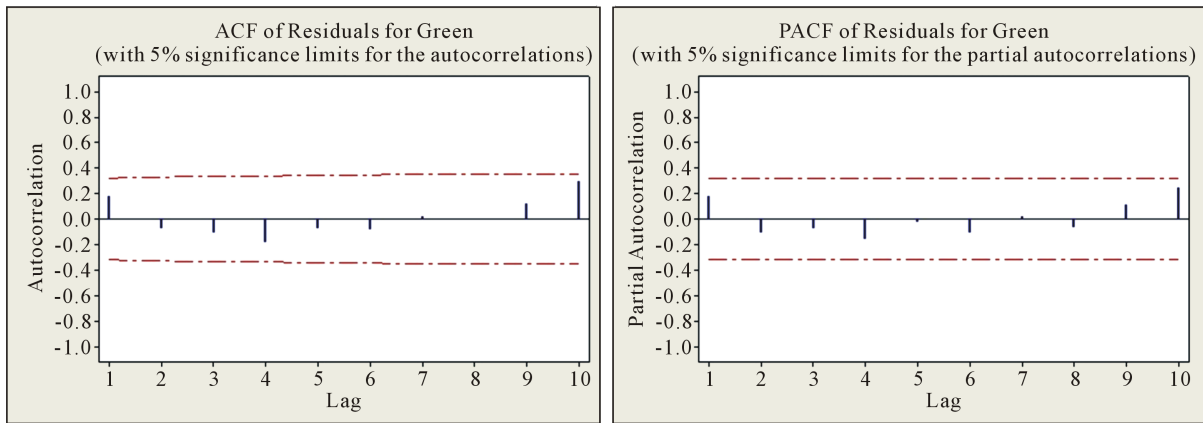


Figure 9. ACF and PACF for residuals of the AR(1) process.

low and diminish as the number of lags increases.

Evidence therefore exists to support that the residuals follow a white noise process and the AR(1) is a robust representation of the observed time series.

The AR(1) process is defined by

$$y_t = 368.1 + 0.694y_{t-1}$$

The forecast for the process will converge onto the process mean, $\mu = 1203$.

The procedure applied to the Yellow shell consumption also resulted in an AR(1) process. Though the details are omitted, the AR(1) process is defined by

$$y_t = 86.73 + 0.8665y_{t-1}$$

The forecast for the process will converge onto the process mean, $\mu = 649$.

5.2. Exponential Smoothing

The application the first order exponential smoothing method requires a choice of smoothing factor λ and the number of observations to smooth against. There is no analytical method to determine the optimum choice of smoothing factor and it is necessary to investigate various levels of λ and choose the smoothing factor that minimises the squared sum of the forecast errors e_t defined by

$$SS_E(\lambda) = \sum_{t=1}^T e_t^2.$$

Similarly there is no method to determine the optimum number of observations to forecast against. The influence of past the observations decays geometrically over time and the impact of the decay can be evaluated against a set of observations.

Table 7 returns generated forecasts for the varying levels of λ against both the Green and Yellow shell data where the full set of 109 observations are used. Due to the magnitude of errors, the square root of $SS_E(\lambda)$ is returned in **Table 7**.

There is a distinct contrast between the two forecasts. Forecast accuracy for the Green consumption increases as λ increases and is a consequence of the volatility of the series. For the Yellow consumption forecast accuracy is assured at $\lambda = 0.4$, where $SS_E(\lambda)$ takes the minimum value. Evident for both time series is that the forecast is influenced by the level of the smoothing factor λ . The choice of smoothing factor for the Yellow forecast is defined by the existence of a minimum $SS_E(\lambda)$. The choice of smoothing factor for the Green forecast is not so clear. By inspection of **Table 6**, choosing $\lambda = 0.6$ would appear appropriate as the forecast reduces at $\lambda = 0.8$ drops slightly.

Table 8 returns the forecast generated at different observation levels for specific choices of smoothing factor λ .

Variation in the forecast level only becomes apparent at the lower number of observations and the variation is

insignificant relative to the level of each forecast and the width of the confidence intervals.

The evidence from **Table 7** and **Table 8** suggests that the level of smoothing factor is critical with respect to generating a robust forecast while the choice of the number of observations to forecast against is not critical.

5.3. Monte Carlo Simulation

Table 9 returns output from the simulation programme for a range of sample periods for both the Green and Yellow shell consumption. What is clear from both sets of simulations is that for any sample period the simulated forecast signals are consistent. For the Green simulation as the sample period increases, volatility of the consumption is exposed to the simulation. As the sample period increases, the standard deviation increases resulting in an inflated forecast relative to the sample mean. For the Yellow simulation, as the sample period is

Table 7. Forecasts generated against smoothing factor level.

Green Forecast Summary (N = 109)						Yellow Forecast Summary (N = 109)					
λ	Forecast	LCL	UCL	Con Int	Sqrt ($SS_E(\lambda)$)	λ	Forecast	LCL	UCL	Con Int	Sqrt ($SS_E(\lambda)$)
0.2	1095.82	-607	2798	3405	9352	0.2	502.18	-246	1251	1497	4273
0.4	1209.07	-218	2636	2855	8048	0.4	644.26	-13	1302	1315	3958
0.6	1265.79	-16	2547	2563	7423	0.6	752.99	107	1399	1292	4034
0.8	1262.91	37	2489	2452	7112	0.8	824.60	166	1483	1318	4256

Table 8. Forecasts generated against number of observations.

Green					Yellow				
$\lambda = 0.6$	Forecast	LCL	UCL	Con Int	$\lambda = 0.4$	Forecast	LCL	UCL	Con Int
N = 109	1265.789	-15.6988	2547.277	2562.975	N = 109	644.2553	-13.0679	1301.579	1314.646
N = 50	1265.789	-98.4743	2630.052	2728.526	N = 50	644.2553	100.0596	1188.451	1088.391
N = 20	1265.789	251.1019	2280.476	2029.374	N = 20	644.2454	171.9441	1116.547	944.6025
N = 10	1265.889	717.1508	1814.628	1097.477	N = 10	643.8595	21.71238	1266.007	1244.294

Table 9. Simulated forecast output.

Green Simulation							
Period	Mean	St Dev	Simulated Forecast Signal				
N = 10	1055.70	271.31	1014	1016	1010	1019	1021
N = 15	933.53	483.71	921	918	924	917	926
N = 20	1013.30	543.61	1271	1271	1269	1277	1264
N = 25	1289.44	907.56	1760	1773	1759	1769	1779
Yellow Simulation							
Period	Mean	St Dev	Simulated Forecast Signal				
N = 10	461.00	279.72	475	478	486	475	481
N = 15	344.40	292.29	442	445	448	454	445
N = 20	328.95	257.37	422	419	415	416	425
N = 25	391.08	303.48	503	498	501	507	496

extended, the standard deviation has not increased significantly reflecting that over extended sample periods the distribution of observations has a greater consistency.

The simulation process is sensitive to excessive volatility resulting in an inflated forecast signal. In principle the affect of the volatility can be overcome by reducing an excessive observation to say within one or two standard deviations of the sample mean. The difficulty in this approach is creating a consistent rule that would apply to all excessive observations. Setting a trigger point at for example 2.5 standard deviations above the sample, an observation at 2.6 standard deviations would be excessively reduced while an observation at 3.5 standard deviations may not be reduced enough.

5.4. Comparison of Forecasting Methods

Table 10 returns the observations of the shell consumption for the following 12 weeks beyond the original observations and the respective graphs of the time series is returned in **Figure 10**.

A summary of the output from each method is presented in **Table 11** where the totals of each forecast are compared to the total of the observed values over the forecast period.

The forecast signals are reasonably close to one another. Though the simulated forecast signal for the Yellow shell is significantly lower than the signals generated by the time series methods, the total forecast is in excess of the observed consumption over the forecast period. The consistency of the forecast signals imply that each modelling processes is robust so indicating the reliability of the signals to generate purchase orders for the shells.

Table 10. Observed shell consumption over forecast period.

	1	2	3	4	5	6	7	8	9	10	11	12	Total
Green	1433	1449	234	305	319	308	362	721	795	877	791	941	8535
Yellow	1115	509	111	491	445	374	387	603	457	432	89	59	5072

Table 11. Summary of forecast output.

Method	Shell	Forecast Signal	Total Forecast	Observed Consumption	Excess
Autoregressive	Green	1203	14,436	8535	5901
	Yellow	649	7788	5072	2716
Exponential Smoothing	Green	1265	15,180	8535	6645
	Yellow	644	7728	5072	2656
Simulation	Green	1275	15,300	8535	6765
	Yellow	475	5700	5072	628

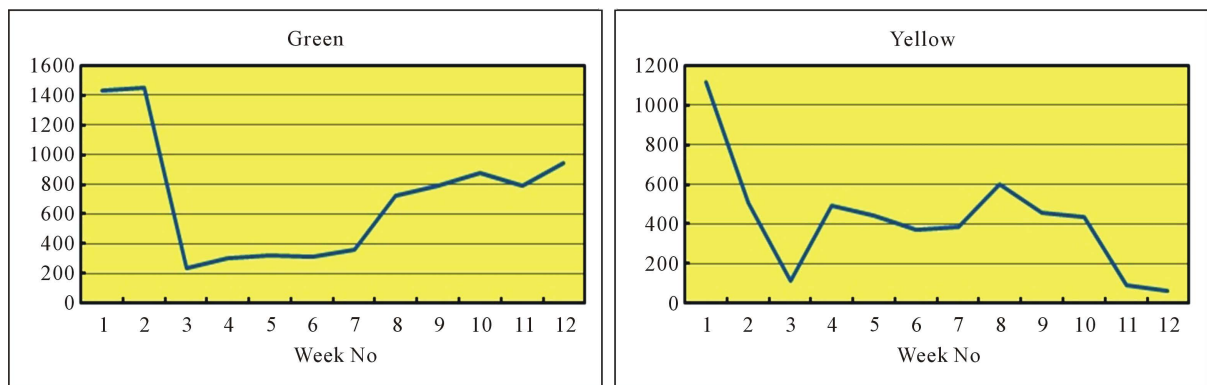


Figure 10. Time series observations over forecast period.

6. Conclusions

Historically, forecasting future demand of the crankshaft shells has proven exceptionally difficult. The purchasing professionals within the case study environment having no experience of applying formal forecasting methods placed orders on the crankshaft shell suppliers that were effectively a “best guess”. To mitigate the need to guess, formal time series analysis was assessed as a suitable approach to modelling demand along with a Monte Carlo simulation method coded in Microsoft Excel VBA.

The forecasts generated by the exponential smoothing and autoregressive process are remarkably close. The simulation process though proving to be sensitive to volatile consumption with judgement of the physical time series, an appropriate forecast signal can be found.

The advantage of the exponential smoothing method lies in its simplicity of execution. The method is open to coding in VBA and would therefore negate the need for specialised software. Though for a volatile time series, it may not be possible to find a smoothing factor that will maximise forecast accuracy. However, choosing a smoothing factor between 0.5 and 0.7 should provide an appropriate forecast signal.

The ARIMA process requires interaction between the user and the modelling process as the ACF and PACF have to be inspected to determine the structure of the model. Though in this study, the autoregressive models were basic AR(1) models, future trajectories may follow more complex ARIMA models. It is recommended that at least 30 observations of a time series are needed to enable the generation of a meaningful forecast. In the current study, there is over two years of data available so the restriction did not apply. However, for the build of a new engine, the application of the ARIMA model would be impeded until there are enough observations to forecast against. Empirical evidence from this study confirms that the exponential smoothing model can produce meaningful forecasts with at least 10 observations and would therefore provide a more robust method of forecasting with limited data sets.

The simulation method provides a seamless process to generate forecast signals and offers additional functionality to generate orders. Execution of the simulation for the complete shell requirement is efficient and completes generally in less than two minutes. The simulation is sensitive to volatile demand and can over inflate the forecast signal.

Monte Carlo methods have proved effective to model a diverse range of complex applications. While the method is consistent, the execution of the method to a specific application has to be tailored to that application. The functionality of the simulation method created within this study is restricted to modelling the forecast consumption of the crankshaft shells. Conversely, the time series methods are universally applicable.

Inspection of **Table 11** confirms that for each of methods, the forecast signals are close to each other and so with respect to forecast accuracy there is no one best method. Moreover, since the forecast signals are close to one another, the purchasing professionals are confident that a robust forecast signal is being generated.

Currently, the appropriate choice of forecasting tool is the simulation method as it provides a seamless process not only to generate the forecast signal but also generate the orders. Building a VBA programme around the exponential smoothing process should in principle provide the functionality provided by the simulation method.

Inventory profiles exhibited by the consumption of the crankshaft shells are rare within the discipline of Operations Management. Conventional though rigorous methods of inventory control that include Materials Requirement Planning and Just-in-Time Kanban systems do not apply to the procurement of the crankshaft shells. Moreover, due to the rarity of such inventory profiles, there does not appear to be any significant research into this unique area of inventory management. The application of the Monte Carlo simulation method and time series analysis begins to close this gap.

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Appendix 1. Backshift and Differencing Operators

Frequently within the time series literature, time series are presented using difference operator denoted by ∇ and the lag or backshift operator B . Kirchgässner *et al.* (2013) succinctly define the properties of both operators. The essential properties of the operators are replicated as follows:

First order differencing is expressed using the difference operator ∇ such that

$$\nabla y_t = y_t - y_{t-1},$$

Second order differencing is expressed as

$$\nabla^2 y_t = \nabla(y_t - y_{t-1}) = y_t - 2y_{t-1} + y_{t-2}$$

The backshift operator B has the effect of “delaying” a time series by one period, such that

$$By_t = y_{t-1}$$

Applying the backshift operator to y_{t-1} , the following is obtained

$$By_{t-1} = B(By_t) = B^2 y_t$$

In general

$$B^k y_t = y_{t-k} \tag{A1}$$

Applying property (A1) to a time series of the form

$$y_t = \varepsilon_t + \theta_1 \varepsilon_{t-1} + \theta_2 \varepsilon_{t-2} + \dots + \theta_q \varepsilon_{t-q}$$

yields a polynomial in B such that

$$y_t = (1 + \theta_1 B + \theta_2 B^2 + \dots + \theta_q B^q) \varepsilon_t = \Theta(B) \varepsilon_t$$

The backshift operator can be related to a first order difference in the following way

$$\nabla y_t = y_t - y_{t-1} = y_t - By_t = (1 - B) y_t$$

Second order differencing

$$\nabla^2 y_t = \nabla(\nabla y_t) = (1 - B)(1 - B) y_t = (1 - B)^2 y_t$$

If a series is differenced d times then it can be shown that

$$\nabla^d y_t = (1 - B)^d y_t$$