# Dealing with Spatial Heterogeneity in Pointwise to Gridded Data Comparisons

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#### 14 Abstract

Most studies on validation of satellite trace gas retrievals or atmospheric chemical transport models 15 16 assume that pointwise measurements, which roughly represent the element of space, should compare well with satellite (model) pixels (grid box). This assumption implies that the field of 17 18 interest must possess a high degree of spatial homogeneity within the pixels (grid box), which may 19 not hold true for species with short atmospheric lifetimes or in the proximity of plumes. Results of 20 this assumption often lead to a perception of a nonphysical discrepancy between data, resulting 21 from different spatial scales, potentially making the comparisons prone to overinterpretation. 22 Semivariogram is a mathematical expression of spatial variability in discrete data. Modeling the semivariogram behavior permits carrying out spatial optimal linear prediction of a random process 23 field using kriging. Kriging can extract the spatial information (variance) pertaining to a specific 24 25 scale, which in turn translating pointwise data to a gridded space with quantified uncertainty such that a grid-to-grid comparison can be made. Here, using both theoretical and real-world 26 27 experiments, we demonstrate that this classical geostatistical approach can be well adapted to 28 solving problems in evaluating model-predicted or satellite-derived atmospheric trace gases. This 29 study suggests that satellite validation procedures using the present method must take kriging 30 variance and satellite spatial response functions into account. We present the comparison of Ozone 31 Monitoring Instrument (OMI) tropospheric NO<sub>2</sub> columns against 11 Pandora Spectrometer Instrument (PSI) systems during the DISCOVER-AO campaign over Houston. The least-squares 32 fit to the paired data shows a low slope (OMI= $0.76 \times PSI+1.18 \times 10^{15}$  molecules cm<sup>-2</sup>, r<sup>2</sup>=0.67) 33 which is indicative of varying biases in OMI. This perceived slope, induced by the problem of 34 35 spatial scale, disappears in the comparison of the convolved kriged PSI and OMI  $(0.96 \times PSI+0.66 \times 10^{15} \text{ molecules cm}^2, r^2=0.72)$  illustrating that OMI possibly has a constant 36 systematic bias over the area. To avoid gross errors in comparisons made between gridded data 37 38 versus pointwise measurements, we argue that the concept of semivariogram (or spatial auto-39 correlation) should be taken into consideration, particularly if the field exhibits a strong degree of 40 spatial heterogeneity at the scale of satellite and/or model footprints.

#### 42 **1. Introduction**

43 Most of the literature on validation of satellite trace gas retrievals or atmospheric chemical transport models assume that geophysical quantities within a satellite pixel or a model grid box 44 45 are spatially homogeneous. Nevertheless, it has long been recognized that this assumption can often be violated; spatially coarse atmospheric models or satellites are often not able to represent 46 features, nor physical processes, transpiring at fine spatial scales. Janjic et al. [2016] used the term 47 48 of representation error to describe this complication. They posit that this problem is a result of 49 two combined factors: unresolved spatial scales and physiochemical processes. To elaborate on this definition, let us assume that an atmospheric model simulating CO<sub>2</sub> concentrations can 50 51 represent the exact physiochemical processes but is fed with a constant CO<sub>2</sub> emission rate. This 52 model obviously cannot resolve the spatial distribution of  $CO_2$  concentration because we use an unresolved emission input. As another example, if we know the exact rates of CO<sub>2</sub> emissions but 53 use a model unable to resolve atmospheric dynamics, the spatial distribution of CO<sub>2</sub> concentrations 54 55 will be unrealistic due to unresolved physical processes.

56 Numerous scientific studies have reported on this matter. The simulations of short lifetime 57 atmospheric compounds such as nitrogen dioxide (NO<sub>2</sub>), isoprene, formaldehyde (HCHO), and 58 the hydroxyl radical (OH) have been found to be strongly sensitive to the model spatial resolution 59 [Vinken et al., 2011; Valin et al., 2011; Yu et al., 2016; Pan et al., 2017]. Likewise, the performance of weather forecast models in resolving non-hydrostatic components heavily relies on both model 60 61 resolution and parametrizations used. For example, when Kendon et al. [2014], Souri et al. [2020a], and Wang et al. [2017] defined a higher spatial resolution in conjunction with more 62 elaborate model physics, they were able to more realistically simulate extreme or local weather 63 64 phenomena such as convection and sea-land breeze circulation.

The spatial representation issue is not only limited to models. Satellite trace gas retrievals 65 optimize the concentration of trace gases and/or atmospheric states to best match the observed 66 67 radiance using an optimizer along with an atmospheric radiative transfer model. This procedure requires various inputs such as surface albedo, cloud and aerosol optical properties, and trace gas 68 profiles, all of which come with different scales and representation errors. Moreover, the radiative 69 transfer model by itself has different layers of complexity with regards to physics. A myriad of 70 71 studies have reported that satellite-derived retrievals underrepresent spatial variability whenever the prognostic inputs used in the retrieval are spatially unresolved [e.g., Russell et al., 2011; 72 73 Laughner et al., 2018; Souri et al., 2016; Goldberg et al., 2019; Zhao et al., 2020]. Additionally, 74 the large footprint of some sensors relative to the scale of spatial variability of species inevitably 75 leads to some degree of the representativity issues [e.g., Souri et al., 2020b, Tang et al., 2021; Judd et al., 2020]. It is because of this reason that several validation studies resorted to downscaling 76 77 their relatively coarse satellite observations using high-resolution chemical transport models so that they could compare them to spatially finer datasets such as in-situ measurements [Kim et al., 78 79 2018; Choi et al., 2020]. Nonetheless, their results largely arise from modeling experiments which 80 might be biased.

The validation of satellites or atmospheric models is widely done against pointwise measurements. Mathematically, a point is an element of space. Hence, it is not meaningful to associate a point with a spatial scale. If one compares a grid box to a point sample (i.e., apples to oranges), they are assuming that the point is the representative of the grid box. At this point, the fundamental question is: can the average of the spatial distribution of the underlying compound be represented by a single value measured at a subgrid location? This question was answered in Matheron [1963]. He advocated the notion of the semivariogram, a mathematical description of

the spatial variability, which finally led to the invention of kriging, the best unbiased linear 88 89 estimator of a random field. A kriging model can estimate a geophysical quantity in a common grid. This is not exclusively special; a simple interpolation method such as the nearest neighbor 90 91 has the same purpose. The power of kriging lies in the fact that it takes the data-driven spatial variability information into account and informs an error associated with the interpolated map. 92 93 This strength not only makes kriging a relatively superior model over simplified interpolation 94 methods, but also reflects the level of confidence pertaining to spatial heterogeneity dictated by 95 both data and the semivariogram model used through its variance [Chilès and Delfiner, 2009].

96 Different studies leveraged this classical geostatistical method to map the concentrations 97 of different atmospheric compounds at very high spatial resolutions [Tadíc et al., 2017; Li et al., 98 2019; Zhan et al., 2018; Wu et al., 2018]; To the best of our knowledge, Swall and Foley [2009] is the only study that used kriging for a chemical transport model validation with respect to surface 99 ozone. They suggested that kriging estimation should be executed in grids rather than discrete 100 points. Kriging uses a semivariogram model in a continuous form. Optimizing the kriging grid size 101 102 (i.e., domain discretization) at which the estimation is performed is an essence to fully obtaining the maximum spatial information from data. Another important caveat with Swall and Foley 103 104 [2009] is that averaging discrete estimates (points) to build grids is not applicable for remote sensing data. Depending on the optics and the geometry, the spatial response function can 105 transform from an ideal box (simple average) to a sophisticated shape such as a super Gaussian 106 107 function (weighted average) [Sun et al., 2018]. Moreover, the footprint of satellites is not spatially constant. We will address these complications in this study using both theoretical and real-world 108 experiments. 109

Our paper is organized with the following sections. Sections 2 is a thorough review of the concept of the semivariogram and kriging. We then provide different theoretical cases, their uncertainty, sensitivities with respect to difference tessellation, grid size, and the number of samples. Section 3 proposes a framework for satellite (model) validation using sparse point measurements and elaborates on the representation error using idealized experiments. Sections 4 introduces several real-world experiments.

#### 116 2. Semivariogram and Ordinary Kriging Estimator

#### 117 2.1. Definition

118 The semivariogram is a mathematical representation of the degree of spatial variability (or 119 similarity) in a function describing a regionalized geophysical quantity (*f*), which is defined as 120 [Matheron, 1963]:

$$\gamma(h) = \frac{1}{2V} \iiint_{V} \left[ f(x+h) - f(x) \right]^2 dV \tag{1}$$

where x is a location in the geometric fields of V, f(x) is the value of a quantity at the location of x, and **h** is the vector of distance. If discrete samples are available rather than the continuous field, the general formula can be simplified to the experimental semivariogram defined as:

$$\gamma(\boldsymbol{h}) = \frac{1}{2N(\boldsymbol{h})} \sum_{|x_i - x_j| - |\boldsymbol{h}| \le \varepsilon} [Z(x_i) - Z(x_j)]^2$$
(2)

124 where  $Z(x_i)$  (and  $Z(x_j)$ ) is discrete observations (or samples), N(h) is the number of paired 125 observations separated by the vector of h. |.| operator indicates the length of a vector. The condition 126 of  $|x_i - x_j| - |h| \le \varepsilon$  is to allow certain tolerance for differences in the length of the vector. For 127 simplicity, we only focus on an isotropic case meaning we rule out the directional (or angular) 128 dependency in  $\gamma(h)$ . Under this condition, the vector of h becomes scalar (h = |h|).

- 129 If a reasonable number of samples is present, one can describe  $\gamma(h)$  through a regression 130 model (e.g., Gaussian or spherical shapes). The degree of freedom for this regression is:
  - dof = N p
- 131 where *p* is the number of parameters defined in the model. For instance, to fit a Gaussian function

(3)

- to the semivariogram with three parameters (p=3), three paired (N=3) observations are required at
- 133 minimum. Different regression models can be used to describe  $\gamma(h)$  depending on the
- 134 characteristic of the quantity of interest. In this study, we will use a stable Gaussian function:

$$\gamma(h) = a(1 - e^{-(\frac{h}{b})^{c_0}}): c_0 = 1.5$$
(4)  
135 where *a* and *b* are fitting parameters. A non-linear least-squares algorithm based on Levenberg-

Marquardt method will be used to estimate the fitting parameters.
The kriging estimator predicts a value of interest over a defined domain using a semivariogram model derived from samples [Chilès and Delfiner, 2009]. The kriging model is defined as [Matheron, 1963]:

$$Z(x) = Y(x) + m(x)$$
(5)

140 where Y(x) is a zero-mean random function, and m(x) is a systematic drift. If we assume 141  $m(x) = a_0$ , the model is called ordinary kriging. Similar to an interpolation problem, the 142 estimation point  $(\hat{Z})$ , is determined by linearly combining *n* number of samples with their weights 143  $(\lambda_i)$ :

$$\hat{Z} = \sum_{j=1}^{n} \lambda_j Z(x_j) + \lambda_0 \tag{6}$$

- 144 where  $\hat{Z}$  is the estimation,  $\lambda_o$  is a constant weight,  $x_j$  is the location of samples, , and  $Z(x_j)$  is point
- 145 data (i.e., samples). The mean squared error of this estimation can be written as

$$E(\hat{Z} - Z_0)^2 = \operatorname{Var}\left(\hat{Z} - Z_0\right) + \left[\lambda_0 + (\sum_{j=1}^n \lambda_j - 1)a_0\right]^2$$
(7)

146 Where  $Z_0$  is point observations  $(Z_o = Z(x_j), j = 1, 2, ..., n)$ , and  $a_0$  is the mean of Z which is 147 unknown. In order to estimate the weights, we are required to minimize Eq.7, but this cannot be 148 done without knowing the exact value of  $a_0$ . A solution is to assume  $\lambda_0 = 0$  and impose the 149 following condition:

$$\sum_{j=1}^{n} \lambda_j = 1 \tag{8}$$

- 150 This condition warrants  $E(\hat{Z} Z_0)$  be zero and removes the need for the knowledge of  $a_o$ .
- 151 Therefore Eq.7 can be written as

$$E(\hat{Z} - Z_0)^2 = \operatorname{Var}\left(\hat{Z} - Z_0\right) = \sum_{j_1=1}^n \sum_{j_2=1}^n \lambda_{j_1} \lambda_{j_2} \gamma_{j_1 j_2} - 2\sum_{j_1=1}^n \lambda_{j_1} \gamma_{j_1 0} + \gamma_{00}$$
(9)

- 152 where  $\gamma_{j1j2}$  is the spatial covariance between the point observations and  $\gamma_{j10}$  is the spatial
- 153 covariance of between the observations and the estimation point. The spatial covariance is modeled
- by a semivariogram. Using the method of Lagrange multiplier and considering the constraint on
- the weights, Eq.9 can be minimized by solving the following problem [Chilès and Delfiner, 2009]:

$$\begin{pmatrix} \lambda_1 \\ \vdots \\ \lambda_n \\ \mu \end{pmatrix} = \begin{pmatrix} \gamma(x_1 - x_1) \cdots \gamma(x_1 - x_n) 1 \\ \vdots & \ddots & \vdots & \vdots \\ \gamma(x_n - x_1) \cdots \gamma(x_n - x_n) 1 \\ 1 & \cdots & 1 & 0 \end{pmatrix}^{-1} \begin{pmatrix} \gamma(x_1 - x_0) \\ \vdots \\ \gamma(x_n - x_0) \\ 1 \end{pmatrix}$$
(10)

where  $\mu$  is the Lagrange parameter and  $x_0$  is the location of estimation. The first term in the right hand side of this equation shows the spatial variability described by the semivariogram model among samples, whereas the second term indicates the modeled variability between samples and the estimation point. The unknowns (the left hand side of the equation) have a unique solution if, and only if, the semivariogram model is positive definite and the samples are unique [Chilès and Delfiner, 2009]. The estimation error can be obtained by

$$\sigma^{2} = E(\hat{Z} - Z_{0})^{2} = \sum_{j=1}^{n} \lambda_{j} \gamma_{j0} - \mu$$
(11)

162 This equation is an important component in the kriging estimator. Not only can we estimate Z(x)163 given a selection of data points, but also an uncertainty associated with such estimation can be 164 provided.

#### 165 2.2. Theoretical Cases

#### 166 *2.2.1. Sensitivity to spatial variability of the field*

The present section illustrates the application of ordinary kriging for several numerical 167 cases. Five idealized cases are simulated in a grid of 100×100 pixels, namely, a constant field (C1), 168 a ramp starting from zero in the lower left to higher values in the upper right (C2), an intersection 169 with concentrated values in four corridors (C3), a Gaussian plume placed in the center (C4), and 170 multiple Gaussian plumes spread over the entire domain (C5). We randomly sample 200 data 171 points from each field as is, and successively create the semivariograms in 100 binned distances. 172 173 Except C1, which lacks a spatial variability thus  $\gamma(h) = 0$ , other semivariograms are fit with the stable Gaussian function. Using the semivariogram model, we optimize Eq.10 to estimate  $\hat{Z}(x)$ 174 175 for each pixel (i.e., 100×100) with the estimation errors based on Eq.11. Figure 1 depicts the truth field (Z(x)), semivariograms made from the samples, estimated values  $(\hat{Z}(x))$ , difference of Z(x)176 and  $\hat{Z}(x)$ , and error associated with the estimation. 177

As for C1, the uniformity results in a constant semivariogram leading the estimation to be identical to the truth. This estimation signifies the unbiased characteristic of ordinary kriging. C1 is never met in reality, however, it is possible to assume some degree of uniformity among data restrained to background values; a typical example of this can be seen in the spatial distribution of a number of trace gases in pristine environments such as NO<sub>2</sub> [e.g., Wang et al., 2020] and HCHO [Wolfe et al., 2019]. Under this condition, any data point within the field (i.e., the satellite footprint) can be assumed to be representative of the spatial variability in truth.

Concerning C2, the semivariogram shows a linear shape meaning data points at larger 185 distances exhibit larger differences. Generally geophysical samples are uncorrelated at large 186 187 distances, thereby one expects the semivarioram to increase more slowly as the distance gets 188 further. The steady increase in  $\gamma(h)$  is indicative of a systematic drift in the data invalidating the assumption of  $m(x) = a_0$ . In many applications, a simple polynomial can explain m(x) and 189 190 subsequently be subtracted from the data points. An example of this problem is tackled by Onn 191 and Zebker [2006]; it concerns the spatial variability of water vapor columns measured by GPS 192 signals. Onn and Zebker [2006] observed a strong relationship between the water vapor columns and GPS altitudes resulting from the vertical distribution of water vapor in the atmosphere. 193 194 Because of this complication, a physical drift model describing the vertical dependency was fit and removed from the measurements so that they could focus on the horizontal fluctuations. In terms of C2, one can effortlessly reproduce Z(x) by fitting a three-dimensional plane to barely three samples, indicating that the semivariogram is of little use.

198 C3 is an example of an extremely inhomogeneous field manifested in the stabilized 199 semivariogram at a value of  $\gamma \sim 500$ , called the sill [Chilès and Delfiner, 2009], indicating 200 insignificant information (variance) from the samples beyond this distance ( $\sim 20$ ), called the range. 201 Range is defined as the separation distance at which the total variance in data is extracted. The 202 smaller the range is, the more heterogeneous the samples will be. While the estimated field roughly captures the shape of the intersections, it is spatially distorted at places with relatively sparse data 203 204 points. The kriging model error is essentially a measure of the density of information. It converges 205 to zero in the samples location and diverges to large values in gaps.

206 C4 is a close example of a point source emitter with faint winds and turbulence. The semivariogram exhibits a bell shape. As samples get further from the source, the variance diverges, 207 208 stabilizes, and then sharply decreases. This is essentially because many data points with low 209 values, apart from each other, have negligible differences. This tendency is recognized as the hole 210 effect which is characterized for high values to be systemically surrounded by low values (and 211 vice versa). It is possible to mask this effect by fitting a semivariogram model stabilizing at certain 212 sill (like the one in Figure 1). Nonetheless, if the semivariogram shows periodic holes, the fitted 213 model should be modified to a periodic cosine model [Pyrcz and Deutsch, 2003].

The last case, C5, shows a less severe case of the hole effect previously observed in C4. This is due to the presence of more structured patterns in different parts of the domain. The range is roughly twice as large as the previous case (C4) denoting that there is more information (variance) among the samples at larger distances. A number of experiments using this particular case will be discussed in the following subsections.

#### 219 *2.2.2. Sensitivity to the number of samples*

220 It is often essential to optimize the number of samples used for kriging. The kriging 221 estimator somewhat recognizes its own capability at capturing the spatial variability through Eq.11. Thus, if the target is spatially too complex and/or the samples are too limited, the estimator 222 essentially informs that  $\hat{Z}(x_0)$  is unreliable through large variance. However, there is a caveat; 223 Y(x) must be a Gaussian random model with a zero mean so that kriging can capture the statistical 224 distribution of  $\hat{Z}$  given the data points. Except this case, the kriging variance can either be 225 underestimated or overestimated depending on the level of skewness of the statistical distribution 226 of Y(x) [Armstrong, 1994]. Figure 2 shows the kriging estimation for C5 using 5, 25, 50, 100, and 227 228 500 random samples in the entire field. Immediately apparent is a better description of the 229 semivariogram when larger number of samples are used, which in turn, results in a better estimation of Z(x). The optimum number of samples to reproduce Z(x) depends on the 230 231 requirement for the relative error  $(\sigma/Z(x))$  being met at a given location.

#### 232 *2.2.3.* Sensitivity to the tessellation of samples

A common application of kriging is to optimize the tessellation of data points for a fixed number of samples to achieve a desired precision. In real-world practices, the objective of such optimization is very purpose-specific, for example, one might prefer a spatial model representing a certain plume in the entire domain. Different ways for data selection exist [e.g., Rennen, 2008], but for simplicity, we focus on four categories: purely random, stratified random, a uniform grid, and an optimized tessellation. Figure 3 demonstrates the estimation of C5 using 25 samples chosen based on those four procedures.

240 Concerning the random selection, the lack of samples over two minor plumes cause the 241 estimation to deviate largely from the truth. While a random selection may seem to be practical because it is independent of the underlying spatial variability, it can suffer from under sampling 242 243 issues, thus being inefficient. As a remedy, it might be advantageous to group the domain into 244 similar zones and randomly sample from each, which is commonly known as stratified random 245 selection. We classify the domain into four zones by running the k-mean algorithm on the 246 magnitudes of Z(x) (not shown) and randomly sample six to seven points from each one (total 25). 247 We achieve a better agreement between the estimated field and the truth because we exploited some prior knowledge (here the contrast between low and high values). 248

As for the uniform grid, we notice that there are fewer data points in the semivariogram stemming from redundant distances which is indicative of correlated information. Nonetheless, if the desired tessellation is neutral with regard to location meaning that all parts of the domain is equal of scientific interest, the uniform grid is the most optimal design for the prediction of Z(x)under an ideally isotropic case. A mathematical proof for this claim can be found in Chilès and Delfiner [2009].

To execute the last experiment, we select 25 random samples for 1000 times and find the optimal estimation by finding the minimum sum of  $|\hat{Z}(x_0) - Z(x)|$ . It is worth mentioning that the optimized tessellation is essentially a local minimum based on 1000 kriging attempts. The optimized location of samples seems to more clustered over areas with large spatial gradients. Not too surprisingly, we observe the smallest discrepancy between the estimation and the truth.

260 A lingering concern over the application of these numerical experiments is that the truth is 261 assumed to be known. The truth is never known, by this means we may never exactly know how well or poorly the kriging estimator is performing. However, it is highly unlikely for some prior 262 263 understandings or expectations of the truth to be absent. If this is the case, which is rare, a uniform 264 grid should be intuitively preferred to deliver the local estimations of average values in uniform blocks. In contrast, if the prior knowledge is articulated by previous site visits, model predictions, 265 266 theoretical experiments, pseudo-observations, or other relevant data, the tessellation needs to be 267 optimized.

268 It is important to recognize that the uncertainties associated with the prior knowledge 269 directly affects the level of confidence in the final answer. Accordingly, the prior knowledge error 270 should ultimately be propagated to the kriging variance. The determination of the prior error is often done pragmatically. For example, if the goal is to design the location of thermometer sites to 271 272 capture surface temperature during heat waves using a yearly averaged map of surface temperature, it would be wise to specify a large error with this specific prior information to play 273 274 down the proposed design. This is primarily because the averaged map underrepresents such an 275 atypical case. A possible extension of this example would be to use a weather forecast model with 276 quantified errors capable of capturing retrospective heat waves. Although a reasonable forecast in the past does not necessarily guarantee a reasonable one in the future, it is rational to assume for 277 the uncertainty with a new tessellation design using the weather model forecast to be lower than 278 279 that of using the averaged map.

A general roadmap for the data tessellation design is shown in Figure 4. As proven in Chilès and Delfiner [2009], if the field is purely isotropic, the uniform grid is the most intuitive sensible choice when the prior information on the spatial variability is lacking. When the prior knowledge with quantified errors is available, an optimum tessellation can be achieved by running a large number of kriging models with suitable  $\gamma(h)$  and picking the one yielding the minimum difference between the prior knowledge and the estimation. The choice of the cost function (here L1 norm) is purpose-specific. For example, if the reconstruction of a major plume was the goal, using aweighted cost function, geared towards capturing the shape of plume, would be more appropriate.

288 2.2.4. Sensitivity to the grid size

A kriging model can estimate a geophysical quantity at a desired location considering the data-driven spatial variability information. Since the kriging model is practically in a continuous form, the desired locations can be anywhere within the field of *V*. A question is whether or not it is necessary to map the data onto a very fine grid. There is a trade-off between the computational cost and the accuracy of the interpolated map. The range of the underlying semivariogram helps in finding the optimal solution. The greater the range (i.e., a more homogeneous field), the less important to map the data in a finer grid.

296 Figure 5a depicts an experiment comparing the estimates of C2 at different grid sizes with 297 the truth. The departure of the estimate from the truth is rather negligible for several coarse grids 298 (e.g.,  $10 \times 10$ ). The homogeneous field, manifested by the large range (Figure 1), allows for a reasonable estimation of Z(x) at coarse resolutions with inexpensive computational costs. Figure 299 300 5b shows the same experiment but on C5 with the optimized tessellation. As opposed to the 301 previous experiment, the estimate substantially diverges from the truth when increasing the grid 302 size, suggesting that a finer resolution should be used for fields with smaller ranges (i.e., 303 heterogeneous fields).

The complexity of directly using the range for choosing the optimal grid size arises from 304 the fact that the level of spatial homogeneity can vary within the domain. In fact, the range is 305 306 derived from a semivariogram model representing a crude estimate of varying ranges occurring at 307 various scales. It is intuitively clear that depending on the degree of heterogeneity, which is 308 spatiotemporally variable, the grid size needs to be adaptively adjusted [Bryan, 1999]. For the sake 309 of simplicity, but at a higher computational cost, we adopt a numerical solution which is to first 310 simulate on a coarse grid, then on a finer one until the difference with respect to the previous grid 311 size across all pixels reaches to an acceptable value (<1%). We name this output (1×1) with the optimized tessellation for C5 as C5opt. 312

313 **3.** Comparison of points to satellite pixels

#### 314 *3.1. Synching the scales between the gridded field and satellite pixels*

To minimize the complications of different spatial scales between two gridded data, we first need to upscale the finer resolution data to match the coarse ones. In case of numerical chemical transport or weather forecast models, the size of the grid box is definitive. Likewise, a satellite footprint, mainly dictated by the sensor design, the geometry, and signal-to-noise requirements [Platt et al., 2021], is known. However, the grid size of the kriging estimation is a variable subject to optimization which has been discussed previously.

321 When we compare the grid size of the kriging estimate to that of a satellite (or a model), three situations arise: First, the kriging spatial resolution is coarser than the satellite, a condition 322 occurring when either the field is homogeneous or the field is under sampled. In situations where 323 324 the field is homogeneous ( $\gamma(h) \cong 0$ ), it is safe to directly compare the data points to the satellite measurements without having to use kriging. If the under sampling is the case (see Figure 2 with 325 5 samples), it is sensible to first investigate if the field is homogeneous within the satellite footprint 326 327 using different data (if any). If the homogeneity is met, we either can compare two datasets without 328 kriging or to match the size of kriging grid cell with the satellite footprint and statistically involve the kriging variance in the comparison (discussed later); nonetheless, the kriging estimate beyond 329 the location of samples must be used with extra caution because their variance very quickly 330 departures from zero to extremely large numbers (see Figure 1). Thus, there is a compromise 331

between increasing the number of paired samples between two datasets and enhancing the level of 332 333 confidence in statistics. If independent observations suggest that there might be large heterogeneity 334 within a satellite footprint, it is strongly advised against quantitatively comparing the points to the 335 satellite observations. Second, the number of samples is fewer than three observations in the field 336 so it is in principal impossible to build a semivariogram. Validating a satellite under this condition is prone to misinterpretation because the spatial heterogeneity cannot be modeled. Nonetheless, if 337 one presumes a good degree of homogeneity within the sensor footprint (such as very high-338 339 resolution remote sensing airborne data), the direct comparison of point measurements might be 340 possible. Third, the satellite footprint is coarser than the kriging estimate. Under this condition, we 341 upscale the kriging map to match the spatial resolution of the satellite using

$$\hat{Z}_c = \hat{Z}_f * S = \int \hat{Z}_f(x) S(x - y) dy$$
(12)

where S is the spatial response function,  $\hat{Z}_c$  is the coarse kriging field, <\*> is the convolution operator, y is shift, and  $\hat{Z}_f$  is the fine field. In discrete form we can rewrite Eq.12 in

$$\hat{Z}_{c}[i,j] = \sum_{m} \sum_{n} \hat{Z}_{f}[i-m,j-n] S[m,n]$$
(13)

where *m* and *n* are the dimension of the response function. The mathematical formulation of S[m, n] for a number of satellites can be represented by two-dimensional super Gaussian functions as discussed in Sun et al. [2018]. Atmospheric models have a uniform response to the simulated values within a grid box, therefore  $S[m, n] = \frac{1}{m \times n} J_{m,n}$ , where *J* is the matrix of ones. In the same way, the kriging variance should be convolved through

$$\sigma_c^2[i,j] = \sum_m \sum_n \sigma_f^2[i-m,j-n] S^2[m,n]$$
(14)

where a superscript of 2 denotes squaring, and  $\sigma_c^2$  and  $\sigma_f^2$  are the kriging variance in the coarse and the fine grids, respectively.

To demonstrate the upscaling procedure, we use C5opt  $(1 \times 1)$  and upscale it at six grid sizes 351 (m,m) of 5×5, 10×10, 15×15, 20×20, 25×25, and 30×30. For simplicity, we consider  $S = \frac{1}{m^2} J_{m,m}$ ; 352 this spatial response function results in averaging the values in the grid boxes. Figure 6 shows the 353 354 resultant map overplotted with the samples along with the error estimation. Two tendencies from 355 this experiment can be identified: First, the discrepancy of the point data and  $\hat{Z}$  is becoming more noticeable as the grid size grows; this directly speaks to the notion of the spatial representativeness; 356 357 large grid boxes are less representative of sub-grid values. Second, the gradients of the field along 358 with the estimation error become smoother primarily due to convolving the field with the spatial response function, which acts as a low pass filter. 359

We further directly compare  $\hat{Z}$  to the samples (i.e., observations) shown in Figure 7. We see an excellent comparison between  $\hat{Z}$  at 1×1 resolution with the observations underscoring the unbiasedness characteristic of the kriging estimator. Conversely, the upscaled field gradually diverges from the observations. This divergence is *the problem of scale*.

#### 364 3.2. Point to pixel vs pixel to pixel

To elaborate on the problem of scale, we design an idealized experiment theoretically validating pseudo satellite observations against some pseudo point measurements. The pseudo satellite observations are created by upscaling the C5 truth (Z) to  $30 \times 30$  grid footprint considering

 $S = \frac{1}{m^2} J_{m,m}$ , meaning that the satellite is observing the truth but in a different scale (Figure S1). 368 The pseudo point measurements are the ones used for C5opt. Figure 8a shows the direct 369 comparison of the satellite pixel with the point observations. By ignoring the fundamental fact that 370 these two datasets are inherently different in nature, displaying the same geophysical quantity by 371 at different scales, we observe a perceived discrepancy ( $r^2=0.64$ ). The comparison suggests a 372 wrong conclusion that the satellite observations are biased-low. This discrepancy is unrelated to 373 374 any observational or physical errors, rendering any physical interpretation of the comparison 375 biased due to spatial-scale differences in the data sets. Figure 8b depicts the comparison of each 376 grid box of the upscaled kriging estimate  $(30 \times 30)$  with that of the satellite. This direct comparison shows a strong degree of agreement ( $r^2=0.98$ ), shaking off the erroneous idea of directly comparing 377 point to gridded data when the field exhibits substantial spatial heterogeneity. 378

379 Yet, the comparison misses an important point: the kriging estimate is considered error-380 free. We attempt to incorporate the kriging variance through a Monte Carlo linear regression 381 method. Here, the goal is to find an optimal linear fit  $(y = ax + b + \varepsilon)$  such that  $\chi^2 =$ 382  $\sum \frac{[y-f(x_i,a,b)]^2}{\sigma_y^2 + a^2 \sigma_x^2}$  is minimized.  $\sigma_y^2$  and  $\sigma_x^2$  are the variances of y (here the satellite) and x (the kriging

variance), respectively. We set the errors of y to zero, and randomly perturb the errors of x based 383 384 on a normal distribution with zero mean and a standard deviation equal to that of kriging estimate 15,000 times. The average of optimized a and b coefficients derived from each fit are then 385 386 estimated and their deviation at 95% confidence interval assuming a Gaussian distribution is 387 determined. Figure 8b,c show the linear fit with and without considering the kriging error estimate. 388 The linear fit without involving the kriging error gives a strong impression that it is nearly perfect, 389 following closely to the paired observations. This is essentially explainable by the primary goal of  $\chi^2$  which is to minimize the L2 norm of residuals  $(y - f(x_i, a, b))$ , portraying a very optimistic 390 picture of the satellite validation. The linear fit considering the kriging errors is different. The 391 uncertainties associated with a and b are larger since x is variable (shown in horizontal error bars). 392 The optimal fit gravitates towards the points with smaller standard deviations as they impose a 393 larger weight. The confidence in the linear fit at higher values is lower due to their errors being 394 395 large. This fit is a more realistic portrayal of the satellite validation.

396 Figure 9 summarizes the general roadmap for satellite (and model) validations against point 397 measurements. To fit the semivariogram with at least two parameters, we are required to have 398 three samples at minimum. Therefore, it is implausible to derive the spatial information from the 399 point data where sampling is extremely sparse (<3 samples within the field). The only case of 400 directly comparing point and satellite pixels is when the field within satellite footprint or the field in general is rather homogeneous confirmed by independent data/models. Having more samples 401 402 allows to acquire some information on the spatial heterogeneity. The information carried by the 403 data is considered more and more robust with increasing the number of samples. Subsequently, the kriging map along with its variance derived from a reasonable semivariogram at an optimized 404 grid resolution should be convolved with the satellite response function so that we can conduct an 405 406 apples-to-apples comparison. A real-world example on the satellite validation will be shown later.

#### 407 4. Real-world experiments

#### 408 4.1. Spatial distribution of NO<sub>2</sub>

We begin with focusing on tropospheric NO<sub>2</sub> columns observed by TROPOMI sensor
[Copernicus Sentinel data processed by ESA and Koninklijk Nederlands Meteorologisch Instituut
(KNMI), 2019; Boersma et al., 2018] at ~13:30 LST. We choose NO<sub>2</sub> primarily due to its spatial
heterogeneity [e.g., Souri et al., 2018; Nowlan et al., 2016, 2018; Valin et al., 2011; Judd et al.,

2020]. We oversample good quality pixels (ga flag>0.75) through a physical-based gridding 413 approach [Sun et al., 2018] over Texas at 3×3 km<sup>2</sup> resolution in four seasons in 2019. We extract 414 samples by uniformly selecting the NO<sub>2</sub> columns in the center of each  $30 \times 30$  km<sup>2</sup> block. The 415 416 semivariogram along with its model are calculated, and then we krige the samples. Figure 10 shows the NO<sub>2</sub> columns map for four different seasons, the semivariogram, the kriging estimates, and the 417 418 differences between the estimate and the field. High levels of NO<sub>2</sub> are confined to cities indicating 419 the sources being predominantly anthropogenic. Wintertime NO<sub>2</sub> columns are larger than 420 summertime mainly due to meteorological conditions and the OH cycle, the major sink of NO<sub>2</sub>. All semivariograms exhibit the hole effect. This is because of high values of NO<sub>2</sub> being 421 422 systematically surrounded by low values. Regardless of the season, we fit the stable Gaussian to variances at distances smaller than 2.5° (~275 km<sup>2</sup>). The  $b_0$  parameter explaining the length scale 423 is found to be 0.94, 0.88, 0.71, and 0.83 degree for DJF, MAM, JJA, and SON, respectively. These 424 numbers strongly coincide with the seasonal lifetime of NO<sub>2</sub> [Shah et al., 2020]; wintertime NO<sub>2</sub> 425 columns are spatially more uniform around the sources thus in relative sense, they are more 426 427 homogeneous (spatially correlated) than those in warmer seasons. On the other hand, the shorter 428 NO<sub>x</sub> lifetime in summer results in a steeper gradient of NO<sub>2</sub> concentrations. This tendency should not be generalized because transport and various NOx sources including biomass burning, soil 429 emissions, and lightning and can have large spatiotemporal variability resulting in different length 430 431 scales in different times of a year. The differences between the kriging estimate and the field show 432 some spatial structures indicating that NO<sub>2</sub> is greatly heterogenous.

#### 433 *4.2. Optimized tessellation over Houston*

The preceding TROPOMI data enabled us to optimize a tessellation of ground-based point 434 435 spectrometers over Houston. Our goal here is to propose an optimized network for winter 2021 given our knowledge on the spatial distribution of NO<sub>2</sub> columns in winter 2019 measured by 436 TROPOMI. The assumption of using a retrospective NO<sub>2</sub> field for informing a hypothetical future 437 campaign is not entirely unrealistic. If we have a consistent number of pixels from TROPOMI 438 between two years, it is unlikely for the spatial variance of NO<sub>2</sub> to be substantially different for 439 440 the same season. We follow the framework proposed in Sect. 2.2.3 involving randomly selecting 441 samples from the field (for 50000 iteration), and calculating kriging estimates for a given number of spectrometers. We then chose the optimum tessellation based on the minimum sum of  $|\hat{Z}(x_0) - \hat{Z}(x_0)|$ 442 443  $Z(x)|_{.}$ 

444 Figure 11 shows the optimized tessellation given 5, 10, 15, and 20 spectrometers over Houston. The Houston plume is better represented with more samples being used. All cases share 445 the same feature; the optimized samples are clustered in the proximity or within the plume. This 446 447 tendency is clearly intuitive. We are required to place the spectrometers in locations where a substantial gradient (variance) in the field is expected. The difference between kriging estimate 448 449 and the TROPOMI observations using 20 samples does not substantially differ in comparison to 450 the one using 15 samples. Therefore, to keep the cost low, a preferable strategy is to keep the number of spectrometers as low as possible while achieving a reasonable accuracy. Based on the 451 452 presented results, the optimized tessellation using 15 samples is preferred among others because it 453 achieves roughly the same accuracy as the one with 20 samples.

## 454 4.3. Validating OMI tropospheric NO<sub>2</sub> columns during DISCOVER-AQ 2013 campaign using 455 Pandora

In order to understand ozone pollution [e.g., Mazzuca et al., 2016; Pan et al., 2017; Pan et al., 2015], characterize anthropogenic emissions [Souri et al., 2016, 2018], and validate satellite data [Choi et al., 2020], an intensive air quality campaign was made in September 2013 over

459 Houston (DISCOVER-AQ). The campaign encompassed a large suite of Pandora spectrometer 460 instrument (PSI) (11 stations) measuring total NO<sub>2</sub> columns with a high precision  $(2.7 \times 10^{14})$ molecules  $cm^{-2}$ ) and a moderate nominal accuracy (2.7×10<sup>15</sup> molecules  $cm^{-2}$ ) under the clear-sky 461 462 condition [Herman et al., 2009]. We remove the observations with an error of >0.05 DU, contaminated by clouds, and averaged them over the month of September at 13:30 LST ( $\pm$  30 463 mins). We attempt to validate OMI tropospheric NO<sub>2</sub> columns version 3.0 [Bucsela et al., 2013] 464 refined in Souri et al. [2016] with the 4-km model profiles. The OMI sensor resolution varies from 465 466  $13 \times 34$  km<sup>2</sup> at nadir to  $\sim 40 \times 160$  km<sup>2</sup> at the edge of the scan line. Biased pixels were removed based on cloud fraction > 0.2, terrain reflectivity > 0.3, and main (xtrack) guality flags =0. Following 467 Sun et al. [2018], we oversample high quality pixels in the month of September 2013 over Houston 468 469 at  $0.2 \times 0.2^{\circ}$  resolution. To remove the stratospheric contributions from PSI measurements, we subtract OMI stratospheric NO<sub>2</sub> ( $2.8\pm0.16 \times 10^{15}$  molecules cm<sup>-2</sup>) from the total columns over the 470 area. Figure 12 shows the monthly-averaged tropospheric NO<sub>2</sub> columns measured by OMI 471 overplotted by 11 PSIs. The elevated NO<sub>2</sub> levels (up to  $\sim 6 \times 10^{15}$  molecules cm<sup>-2</sup>) are seen over the 472 473 center of Houston.

474 We then follow the validation framework shown in Figure 9 in which the number of point 475 measurements and the level of heterogeneity are the main factors in deciding if we should directly 476 compare them to the satellite pixels. Figure 13 shows the monthly-averaged PSI measurements 477 along with the semivariogram and resulting kriging estimate at an optimized resolution ( $\sim 2 \text{ km}^2 =$ 478 13800 data over the entire region) and errors. The distribution of semivariogram suggests that there is a strong degree of spatial heterogeneity, necessitating the use of kriging. We fit a stable Gaussian 479 to the semivariogram resulting in 2.23 ×  $(1 - e^{-(\frac{h}{0.19})^{1.5}})$ . The spatial information (variance) levels 480 off at 0.19° (~21 km) with a maximum variance equal to 2.23 molecules<sup>2</sup> cm<sup>-4</sup>. The measurements 481 beyond this range (21 km) have a minimal weight due to this length scale. It is because of this 482 reason that we see the kriging estimate converges to a fixed value at places being further than this 483 range. The kriging errors of those grid boxes are constantly large (40% relative error). The 484 485 optimum grid size for kriging is found to be  $2 \text{ km}^2$  (<1% difference across all grid boxes). 486 Subsequently, we use the super Gaussian spatial response function described in Sun et al. [2018] to convolve both the kriging estimate and error within (see Figure S2). Figure 14 shows the 487 differences between the kriging estimate and error before and after convolution. The response 488 function (OMI pixel) tends to be on average coarser than 2 km<sup>2</sup> resulting in smoothing of both the 489 490 kriging estimate and error.

491 We ultimately conduct two different sets of comparison: directly comparing PSI to OMI pixels, and comparing convolved kriged PSI to OMI. It is worth noting that PSI measurements are 492 493 monthly-averaged; similarly OMI data are oversampled in a monthly basis. In terms of the PSI, we only account for grid boxes whose kriging error is below  $1.2 \times 10^{15}$  molecules cm<sup>-2</sup> (1193) 494 samples, 8% of total kriging grid boxes). As for the grid-to-grid comparison, the kriging variance 495 is considered in the linear polynomial fitted to the data through the Monte Carlo of chi-square with 496 5,000 iterations. The variability with the OMI stratospheric NO<sub>2</sub> columns ( $0.16 \times 10^{15}$  molecules 497 cm<sup>-2</sup>) is added to the PSI error for both analyses. The left and right panels of Figure 15 show the 498 499 comparisons. As for the direct comparison of actual points (PSI) to pixels (OMI), the PSI measurements indicate a deviation of the slope ( $r^2=0.66$ ) from the unity line. This suggests that 500 501 there is an unresolved magnitude-dependent systematic error. The grid-to-grid comparison not 502 only offers a clearer picture of the distribution of data points, but also it hints at the offset being rather constant ( $0.66\pm0.18\times10^{15}$  molecules cm<sup>-2</sup>; r<sup>2</sup>=0.72). We also observe that the statistics 503 between the satellite and the benchmark are moderately improved. This comparison in general 504

provides an important implication: the varying offsets in a plume shape environment (high to low values) are not necessarily due to variable offsets in the satellite retrieval, as the kriging estimate
 suggests that those varying offsets in point-to-pixel comparison, manifested in slope = 0.76, are a

508 result of varying spatial scales.

#### 509 Summary

510 There needs to be increased attention to the spatial representativity in the validation of satellite (model) against pointwise measurements. A point is the element of space, whereas satellite 511 512 (model) pixels (grid box) are (at best) the product of the integration of infinitesimal points and a normalized spatial response function. If the spatial response function is assumed to be an ideal 513 514 box, the resulting grid box will represent the average. Essentially, no justifiable theory exists to 515 accept that the averaged value of a population should absolutely match with a sample, unless all samples are identical (i.e., a spatially homogeneous field). This glaring fact is often overlooked in 516 the atmospheric science community. At a conceptual level, we are required to translate pointwise 517 data to the grid format (i.e., rasterization). This can be done by modeling the spatial autocorrelation 518 (or semivariogram) extracted from the spatial variance (information) among measured sample 519 520 points. Assuming that the underlying field is a random function with an unknown mean, the best linear unbiased predictions of the field can be achieved by kriging using the modeled 521 522 semivariograms.

523 In this study, we discussed methods for the kriging estimation of several idealized cases. 524 Several key tendencies were observed through this experiment: first, the range corresponded to the 525 degree of spatial heterogeneity; a larger range indicated the less presence of heterogeneity. Second, 526 the kriging variance explaining the density of information guickly diverged from zero to large 527 values when the field exhibited large spatial heterogeneity. This tendency mandates increasing the number of samples (observations) for those cases. Third, while the semivariogram models were 528 constructed from discrete pair of samples, they are mathematically in a continuous form. It is 529 530 because of this reason that we determined the optimal spatial resolution of the kriging estimate by incrementally making the grids finer and finer until a desired precision (=1%) was met. 531

The present study applied kriging to achieve an optimum tessellation given a certain number of samples such that the difference between our prior knowledge of the field, articulated by previous observations, models or theory, and the estimation is minimal. Usually there is uncertainty about the prior knowledge that should be propagated to the final estimates. The optimum tessellation for a range of idealized and real-world data consistently voted for placing more samples in areas where the gradients in the measurements were significant such as those close to point emitters.

539 This study also revisited the spatial representativity issue; it limits the realistic 540 determination of biases associated with satellites (models). In one experiment, we convolved the kriging estimate for a multi-plume field with a box filter but various sizes. The perfect agreement 541 (r=1.0) between the samples (point) and kriging output (pixel) seen at a high spatial resolution 542 543 gradually vanished with coarsening of the resolution of grid boxes (r=0.8). We also directly compared samples (point) with pseudo satellite observations (showing the truth) with a coarse 544 spatial resolution which led to a flawed conclusion about the satellite being biased-low. We 545 546 modeled the semivariogram of those samples, estimated the field using kriging, and convolved with the pseudo-satellite spatial response function. The direct comparison of this output with that 547 of the satellite showed a completely different story suggesting that the data were rather free of any 548 549 bias. A serious caveat with using a spatial model (here kriging) is that it consists of errors: the estimations being further from samples are less certain. It is widely known that discounting the 550

measurement/model errors in true straight-line relationship between data can introduce artifacts.
 To consider the kriging variance in the comparisons we employed a Monte Carlo method on chi square optimization which ultimately allowed us to not only provide a set of solutions within the
 range of the uncertainty of the kriging model, but also to assign smaller weights on gross estimates.

We further validated monthly-averaged Ozone Monitoring Instrument (OMI) tropospheric 555 556 NO<sub>2</sub> columns using 11 Pandora Spectrometer Instrument (PSI) observations over Houston during NASA's DISCOVER-AQ campaign. A pixel-to-point comparison between two dataset suggested 557 558 varying biases in OMI manifested in a slope far from the identity line. By contrast, the kriging estimate from the PSI measurements, convolved with the OMI spatial response function, resulted 559 560 in an inter-comparison slope close to the unity line. This suggested that there was only a constant systematic bias  $(0.66\pm0.18\times10^{15} \text{ molecules cm}^{-2})$  associated with the OMI observations which 561 does not vary with tropospheric NO<sub>2</sub> column magnitudes. 562

The central tenants of satellite and model validation are pointwise measurements. Our experiments paved the way for a clear roadmap explaining how to transform these pointwise datasets to a comparable spatial scale relative to satellite (model) footprints. It is no longer necessary to ignore *the problem of scale*. The validation against point measurements can be carefully conducted in the following steps:

- i. Construct the experimental semivariogram if the number of point measurements allows (usually >= 3 within the field; the field can vary depending on the length scale of the compound).
- 572 ii. Drop the quantitative assessment if the number of point measurements are
  573 insufficient to gain spatial variance and the prior knowledge suggests a high
  574 likelihood of spatial heterogeneity within the field.
- 575 iii. Choose an appropriate function to model the semivariogram.
- 576iv.Estimate the field with kriging (or any other spatial estimator capable of digesting577the semivariogram) and calculate the variance.
- 578 v. Estimate the optimum grid resolution of the estimate.
- vi. Convolve the kriging estimate and its variance with the satellite (model) spatial
  response function (which is sensor specific).
- 581vii.Conduct the direct comparison of the convolved kriged output and the satellite582(model) considering their errors through a Monte Carlo (or a weighted least-squares583method).
- 584

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Recent advances in satellite trace gas retrievals and atmospheric models have helped extend our understanding of atmospheric chemistry but an important task before us in improving our knowledge on atmospheric composition is to embrace the semivariogram (or spatial autocorrelation) notion when it comes to validating satellites/models using pointwise measurements, so that we can have more robust quantitative applications of the data and models.

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#### 597 Author contributions

598 AHS designed the research, executed the experiments, analyzed the data, made all figures, and 599 wrote the paper. KS implemented the oversampling method, provided the spatial response 600 functions, and oversampled TROPOMI data. KC, XL, and MSJ helped with the conceptualization 601 of the study and the interpretation of the results. All authors contributed to discussions and edited 602 the paper.

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788 **Figure 1**. (first column) Five theoretical fields randomly sampled with 200 points (dots), namely, a constant field (C1), a ramp starting from zero in the lower left to higher values in the upper right 789 790 (C2), an intersection with concentrated values in four corridors (C3), a Gaussian plume placed in 791 the center (C4), and multiple Gaussian plumes spread over the entire domain (C5). (second column) 792 the corresponding isotropic semivariograms computed based on Eq.2; the red line shows the stable Gaussian fitted to the semivariogram based on Levenberg-Marquardt method. (third column) The 793 kriging estimate at the same resolution of the truth (i.e.,  $1 \times 1$ ) based on Eq.6. (fourth column) The 794 795 difference between the estimate and the truth. (fifth column) the kriging standard error based on 796 Eq.11.



Figure 2. (first column) The multi-plume case (C5) randomly sampled with different number of
samples (5, 25, 50, 100, and 500), (second column) the corresponding isotropic semivariogram,
(third column) the kriging estimate, (fourth column) the difference between the estimate and the
truth, and (fifth column) the kriging standard error.



Figure 3. The multi-plume case (C5) randomly sampled by four different sampling strategies
using a constant number of samples (25). The sampling strategies include purely random (first
row), stratified random (second row), uniform grids (third row), and an optimized tessellation
proposed based on kriging (fourth row). Columns represent the truth, the isotropic semivariogram,
the kriging estimate, the difference between the estimate and the truth, and the kriging standard
error.



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  815 Figure 4. A schematic illustrating a framework for optimum sampling (tessellation) strategy. The
- 816 prior knowledge refers to any data being able of describing our quantity of interest including site-
- 817 visits, theoretical models, satellite observations, emissions, and etc.
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823 different grid resolutions ranging from  $25 \times 25$  pixel to  $0.2 \times 0.2$ . (b) The kriging estimates of the

824 multi-plume (C5) with optimized samples shown in Figure 3 for different grid resolutions. C2 is

825 more homogeneous than C5, as a result, it is less sensitive to the resolution of the kriging

826 estimate. The optimum grid resolution for C2 is  $10 \times 10$ , whereas it is  $1 \times 1$  for C5. These numbers

are based on observing negligible difference (<1%) between the kriging estimate at the optimum resolution and the one computed at a finer resolution step. We call the optimum output for C5 as C5opt.

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**Figure 6.** (first row) C5Opt outputs convolved with an ideal box kernel with different sizes  $(1 \times 1$  up to  $30 \times 30$ ) overlaid by the C5Opt optimum samples. (second row) the associated kriging errors

convolved with the same kernel. The coarser the resolution is, the larger the discrepancy between
the samples and the estimates is.



**Figure 7.** Illustrating the problem of spatial scale: comparisons of the kriging estimates at seven

- 844 different spatial scales with the samples used for the C5opt estimation. The perceived
- 845 discrepancies are purely due to the spatial representativeness.





Figure 8. (a) the direct comparison of pseudo observations of a satellite observing the C5 case at 852 30×30 resolution versus the 25 samples used for C5opt. (b) same for y-axis, but the point samples 853 854 are transformed to grid boxes using kriging convolved with the satellite spatial response function (ideal box with 30×30 kernel size). The differences in statistics between these two experiments 855 856 speak to the problem of scale. (b) ignores the kriging errors but (c) incorporates them using a Monte Carlo method. Note that the best linear fit has changed indicating that the consideration of 857 the kriging variance is critical. MB = mean bias (point minus satellite), MAB = mean absolute 858 bias, RMSE = root mean square error,  $R^2 = coefficient$  of determination. 859 860



Figure 9. The proposed roadmap for transforming pointwise measurements to gridded data insatellite (model) validation.



8 866 Figure 10. (first column) The spatial distribution of TROPOMI tropospheric NO<sub>2</sub> columns oversampled in four different seasons at  $3 \times 3$  km<sup>2</sup> spatial resolution. (second column) The 867 corresponding semivariogram from samples selected from uniform 30×30 km<sup>2</sup> blocks (shown 868 869 with black dots in the first column) along the fitted stable Gaussian model (red line). (third column) the kriging estimates, and (fourth column) their differences with respect to the 870 871 observations.

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Figure 11. Finding an optimum sample tessellation for wintertime over Houston given different
number of spectrometers (5, 10, 15, and 20).



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**Figure 12.** The spatial distribution of OMI tropospheric NO<sub>2</sub> columns oversampled at the resolution at  $0.2 \times 0.2^{\circ}$  over Houston in September 2013. The plot is overlaid by surface Pandora spectrometer instrument averaged over the same month. The surface measurements originally measured the total columns, therefore we subtract the stratospheric columns provided by the OMI data ( $2.8\pm0.16 \times 10^{15}$ molecules cm<sup>-2</sup>) from the total columns to focus on the tropospheric part.



Figure 13. The Pandora tropospheric NO<sub>2</sub> measurements (made from subtracting the total columns
from the OMI stratospheric NO<sub>2</sub> columns) during September 2013, the corresponding
semivariogram, the kriging estimates, and the kriging standard errors. Note that the semivariogram
suggests a large degree of spatial heterogeneity occurring at different spatial scales.



Figure 14. Convolving both kriging estimates and errors with the OMI spatial response function
formulated in Sun et al. [2018]. The differences against the pre-convolved fields are also depicted.





**902 Figure 15.** (left): the direct comparison of OMI tropospheric NO<sub>2</sub> columns with 11 pointwise 903 Pandora measurements in September 2013 over Houston. (right) same for y-axis, but the PSI 904 measurements are translated to grid boxes using kriging convolved with the OMI spatial response 905 function. PSI tropospheric NO<sub>2</sub> columns are estimated based on subtracting the OMI stratospheric 906 NO<sub>2</sub> columns ( $2.8\pm0.16 \times 10^{15}$  molecules cm<sup>-2</sup>) from the total columns. We only consider kriging 907 estimates whose errors are below  $1.2\times10^{15}$  molecules cm<sup>-2</sup>. The kriging variance is also considered 908 using the Monte Carlo method applied on  $\chi^2$ . The slope has improved after considering the 909 modeled spatial representativeness. MB = mean bias (OMI vs Pandora), MAB = mean absolute 910 bias, RMSE = root mean square error.