

Simple and honest confidence intervals in nonparametric regression

TIMOTHY B. ARMSTRONG

Department of Economics, Yale University

MICHAL KOLESÁR

Department of Economics, Princeton University

We consider the problem of constructing honest confidence intervals (CIs) for a scalar parameter of interest, such as the regression discontinuity parameter, in nonparametric regression based on kernel or local polynomial estimators. To ensure that our CIs are honest, we use critical values that take into account the possible bias of the estimator upon which the CIs are based. We show that this approach leads to CIs that are more efficient than conventional CIs that achieve coverage by undersmoothing or subtracting an estimate of the bias. We give sharp efficiency bounds of using different kernels, and derive the optimal bandwidth for constructing honest CIs. We show that using the bandwidth that minimizes the maximum mean-squared error results in CIs that are nearly efficient and that in this case, the critical value depends only on the rate of convergence. For the common case in which the rate of convergence is $n^{-2/5}$, the appropriate critical value for 95% CIs is 2.18, rather than the usual 1.96 critical value. We illustrate our results in a Monte Carlo analysis and an empirical application.

KEYWORDS. Confidence intervals, regression discontinuity, nonparametric regression.

JEL CLASSIFICATION. C14, C21.

1. INTRODUCTION

This paper considers the problem of constructing confidence intervals (CIs) for a scalar parameter $T(f)$ of a function f , which can be a conditional mean or a density. The scalar parameter may correspond, for example, to a conditional mean, or its derivatives at a point, the regression discontinuity or the regression kink parameter, or the value of a density or its derivatives at a point. A popular approach to estimation of $T(f)$ is to use kernel or local polynomial estimators. These estimators are both simple to implement,

Timothy B. Armstrong: timothy.armstrong@yale.edu

Michal Kolesár: mcolesar@princeton.edu

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and highly efficient in terms of their mean squared error (MSE) properties (Fan (1993), Cheng, Fan, and Marron (1997)). CIs are typically formed by undersmoothing (choosing the bandwidth to shrink more quickly than the MSE optimal bandwidth) or bias-correction (subtracting an estimate of the estimator's bias).

In this paper, we propose a simple alternative approach to forming CIs based on these estimators that is more efficient than both undersmoothing and bias-correction in the sense that it leads to shorter CIs while maintaining coverage over the same parameter space \mathcal{F} for f (which typically places bounds on derivatives of f). In particular, one simply adds and subtracts the estimator's standard error times a critical value that is larger than the usual normal quantile $z_{1-\alpha/2}$, and takes into account the possible bias of the estimator.¹ Asymptotically, these CIs correspond to fixed-length CIs as defined in Donoho (1994), and so we refer to them as fixed-length CIs. We show that the critical value depends only on (1) the order of the derivative that one bounds to define the parameter space \mathcal{F} ; and (2) the criterion used to choose the bandwidth. In particular, if the MSE optimal bandwidth is used with a local linear estimator, computing our CI at the 95% coverage level amounts to replacing the usual critical value $z_{0.975} = 1.96$ with 2.18.

When the criterion for bandwidth choice is the length of the resulting CI, we show that the resulting bandwidth is in fact *larger* than the MSE optimal bandwidth. This contrasts with the work of Hall (1992) and Calonico, Cattaneo, and Farrell (2018) on optimality of undersmoothing. Importantly, these papers restrict attention to CIs that use the usual critical value $z_{1-\alpha/2}$. It then becomes necessary to choose a small enough bandwidth so that the bias is asymptotically negligible relative to the standard error, since this is the only way to achieve correct coverage. Our results imply that rather than choosing a smaller bandwidth, it is better to use a larger critical value that takes into account the potential bias; this also ensures correct coverage regardless of the bandwidth sequence. While the fixed-length CIs shrink at the optimal rate, undersmoothed CIs shrink more slowly. We also show that under smoothness assumptions needed to implement bias-correction, our CIs shrink at a faster rate than bias-corrected CIs, once the standard error is adjusted to take into account the variability of the bias estimate (Calonico, Cattaneo, and Titiunik (2014) show that doing so is important for maintaining coverage). The over-smoothing relative to the MSE optimal bandwidth is relatively modest: under a range of conditions most commonly used in practice, a fixed-length CI centered at the MSE optimal bandwidth is 99% efficient relative to using the CI optimal bandwidth. Therefore, a practically attractive implementation of our CIs is to simply center them around an estimator with MSE optimal bandwidth, rather than reoptimizing the bandwidth for length and coverage of the CI.

A key requirement that underlies our results is the notion of honesty: as in Li (1989), we require that the CIs cover the true parameter asymptotically at the nominal level uniformly over the parameter space \mathcal{F} . Furthermore, we allow this parameter space to grow with the sample size. The notion of honesty is closely related to the use of the minimax

¹An R package implementing our CIs in regression discontinuity designs is available at <https://github.com/kolesarm/RDHonest>.

criterion used to derive the MSE efficiency results: in both cases, one requires good performance uniformly over the parameter space \mathcal{F} . The requirement that the CIs be honest is necessary for good finite-sample performance. In contrast, approaches to inference based on pointwise-in- f asymptotics, such as using bandwidths that optimize the pointwise-in- f asymptotic MSE can lead to arbitrarily poor finite-sample behavior, as we discuss further in Section 4.1. To illustrate the practical importance of this point, we conduct a Monte Carlo study in which we show that commonly used CIs based on plug-in bandwidths that attempt to estimate this pointwise-in- f optimal bandwidth exhibit severe undercoverage, even when combined with undersmoothing or bias-correction.

When the parameter space places a bound M on a derivative of f , our CIs require this bound to be specified explicitly. While this may appear to be a disadvantage of our particular approach, due to impossibility results of Low (1997), Cai and Low (2004), and Armstrong and Kolesár (2018a), this cannot be avoided, regardless of how one forms the CI, without making further restrictions on the function f . In particular, these papers show that, without additional assumptions on the parameter space, one cannot use a data-driven method to estimate M and maintain coverage over the whole parameter space—any other method that appears to avoid making this choice must do so *implicitly*. For example, an apparent advantage of undersmoothing is that it leads to correct coverage for any fixed smoothness constant M . However, as we discuss in detail in Section 4.2, a more accurate description of undersmoothing is that for each sample size n , it implicitly chooses a constant M_n under which coverage is controlled. Given a sequence of undersmoothed bandwidths, we show how M_n can be calculated explicitly. One can then obtain a shorter CI with the same coverage properties by computing a fixed-length CI for the corresponding M_n . Regardless of how one chooses M , the fixed-length CIs we propose are more efficient than undersmoothed or bias-corrected CIs that use the same (implicit or explicit) choice of M . In fact, it follows from the calculations in Donoho (1994) and Armstrong and Kolesár (2018a) that our CIs, when constructed using a length-optimal or MSE-optimal bandwidth, are highly efficient among *all* honest CIs: no other approach to inference can substantively improve on their length, while still maintaining coverage.

As an alternative to choosing M a priori, one can place additional conditions on the function f that allow for an upper bound on M to be estimated. To maintain efficiency of the resulting CI, however, care must be taken in doing so: if M is a bound on the p th derivative, and one imposes a bound \tilde{M} on the $(p+1)$ th derivative in order to estimate M , then the optimal CI will be based on a different estimator and will depend on the new bound \tilde{M} . To avoid such issues, we propose a regularity class that relates a global polynomial approximation to smoothness of the function f near the point of interest, and we show formally that, for this class, one can obtain a valid and highly efficient CI using a global polynomial rule of thumb suggested by Fan and Gijbels (1996). However, given the additional assumptions required by this (or any) data driven choice of M , we recommend that this approach be used as a starting point for sensitivity analysis allowing for other choices of M .

Another approach to data-driven choices of M is to use “self-similarity” conditions, as suggested by Giné and Nickl (2010), which relate the maximum and minimum bias

at different bandwidths. Bull (2012) and Chernozhukov, Chetverikov, and Kato (2014) have obtained rate optimal confidence bands under such conditions, which, like the CIs considered here, use a critical value based on an upper bound on the bias. While these results for confidence bands could be extended to cover the problem of constructing CIs for a scalar parameter, obtaining sharp critical values appears to be very difficult. Indeed, the results of Armstrong (2019) show that the sharp form of such CIs must depend to first order on auxiliary constants used to define self-similarity. Nonetheless, our approach of bounding local smoothness using a global polynomial approximation is inspired by the self-similarity approach taken by this literature, and we see it as being in the same spirit. Schennach (2015) also used an upper bound on the bias based on an estimated smoothness constant. While the coverage of the resulting CIs is pointwise-in- f , it is plausible that the CIs are honest under additional auxiliary conditions, similar in spirit to self-similarity.

In addition to calculating the relative efficiency of CIs constructed using different bandwidths, our results allow us to calculate the relative efficiency of CIs constructed using different kernels. In particular, we show that the relative efficiency of kernels for the CIs we propose is *the same* as the relative efficiency of the estimates in terms of MSE. Thus, relative efficiency calculations for MSE, such as the ones in Fan (1993), Cheng, Fan, and Marron (1997), and Fan, Gasser, Gijbels, Brockmann, and Engel (1997) for estimation of a nonparametric mean at a point (estimation of $f(x_0)$ for some x_0) that motivate much of empirical practice in the applied regression discontinuity literature, translate directly to CI construction. Despite their importance in motivating empirical practice, however, such results are subject to a technical critique about how the parameter space is specified: rather than placing a bound on a derivative of f (a Hölder condition), currently available relative efficiency results place assumptions directly on the error of a Taylor approximation at a particular point, so that some “nonsmooth” functions are in fact not ruled out.² To address this, we derive the minimax performance of local polynomial estimators under Hölder restrictions on f . These results confirm that the local polynomial estimators used in empirical practice are also highly efficient under Hölder restrictions on f . Furthermore, while we focus on asymptotic CIs and relative efficiency, these results include a derivation of the finite-sample worst-case bias of local polynomial estimators under Hölder restrictions, which was used by Kolesár and Rothe (2018) to form finite-sample valid CIs in a fixed-design regression setting. These findings may be of independent interest.

The requirement of honesty is also important to ensure that our concept of optimality is well-defined and consistent. As discussed above, it allows us to consider bandwidth or kernel efficiency for constructing CIs. In addition, it also allows us to formally show that using local polynomial regression of an order that is too high given the amount of smoothness imposed is suboptimal. In contrast, under pointwise-in- f asymptotics, high-order local polynomial estimates are superefficient at every point in the parameter space (see Chapter 1.2.4 in Tsybakov (2009), and Brown, Low, and Zhao (1997)).

²See Imbens and Wager (2019), as well as our discussion in Section 3.2.1 for an elaboration of this critique.

To illustrate the implementation of the honest CIs, we reanalyze the data from [Ludwig and Miller \(2007\)](#), who, using a regression discontinuity design, find a large and significant effect of receiving technical assistance to apply for Head Start funding on child mortality at a county level. However, this result is based on CIs that ignore the possible bias of the local linear estimator around which they are built, and an ad hoc bandwidth choice. We find that, if one bounds the second derivative globally by a constant M using a Hölder class, the uncertainty associated with the effect size is much larger than originally reported, unless one is very optimistic about the constant M , allowing f to only be linear or nearly-linear.

Our results build on the literature on estimation of linear functionals in normal models with convex parameter spaces, as developed by [Donoho \(1994\)](#), [Ibragimov and Khas'minskii \(1985\)](#) and many others. As with the results in that literature, our setup gives asymptotic results for problems that are asymptotically equivalent to the Gaussian white noise model, including nonparametric regression ([Brown and Low \(1996\)](#)) and density estimation ([Nussbaum \(1996\)](#)). Our main results build on the “renormalization heuristics” of [Donoho and Low \(1992\)](#), who show that many nonparametric estimation problems have renormalization properties that allow easy computation of minimax MSE optimal kernels and rates of convergence. Our results hold under essentially the same conditions, which apply in many classical nonparametric settings.

The CIs we consider in this paper are applications of the fixed-length CIs proposed in the context of inference on linear functionals $T(f)$ in Gaussian nonparametric regression by [Donoho \(1994\)](#), which have also been studied recently in [Armstrong and Kolesár \(2018a\)](#), and in contemporaneous and subsequent work by [Kolesár and Rothe \(2018\)](#) and [Imbens and Wager \(2019\)](#). In contrast to the finite-sample approach taken in these papers, we focus on asymptotic results, and we also allow $T(f)$ to be nonlinear. Instead of imposing the nonparametric regression model, we require a renormalization condition (see equation (4) below) that allows us to apply the “renormalization heuristics” of [Donoho and Low \(1992\)](#); we are thus able to cover settings such as density estimation or estimation of a bidder valuation in first-price auctions (see Appendix C in the Online Supplemental Material). Our asymptotic approach allows for simplifications that deliver our main relative efficiency results. These efficiency results are different from and complementary to the asymptotic form of the efficiency bounds given in [Donoho \(1994\)](#) and [Armstrong and Kolesár \(2018a\)](#): whereas we consider relative efficiency of estimators and fixed-length CIs based on different kernels and bandwidths, [Donoho \(1994\)](#) and [Armstrong and Kolesár \(2018a\)](#) bound the scope for efficiency gains from CIs that do not fall into this class. [Donoho \(1994\)](#) and [Armstrong and Kolesár \(2018a\)](#) found that the scope for further improvement is small, which motivates our focus on this class of estimators and CIs. See Remark 2.3 for further discussion.

The rest of this paper is organized as follows. Section 2 gives the main results. Section 3 applies our results to inference at a point, sharp and fuzzy RD, and it discusses practical implementation issues, including a rule of thumb for choosing M . Section 4 gives a theoretical comparison of our fixed-length CIs to other approaches, and Section 5 compares them in a Monte Carlo study. Finally, Section 6 presents an empirical application based on [Ludwig and Miller \(2007\)](#). Appendix A gives proofs of the results in

Section 2. Additional results are collected in the Appendices in the Online Supplemental Material (Armstrong and Kolesár (2020)).

2. GENERAL RESULTS

We are interested in a scalar parameter $T(f)$ of a function f , which is typically a conditional mean or a density. The function f is assumed to lie in a function class $\mathcal{F} = \mathcal{F}(M)$, which places “smoothness” conditions on f , where M indexes the level of smoothness. We focus on classical nonparametric function classes, in which M corresponds to a bound on a derivative of f of a given order. We allow $M = M_n$ to grow with the sample size n .

We have available a class of estimators $\hat{T}(h; k)$, indexed by a bandwidth $h = h_n > 0$ and a kernel k . Let $\widehat{\text{se}}(h; k)$ denote the standard error of $\hat{T}(h; k)$, an estimate of its standard deviation $\text{sd}_f(\hat{T}(h; k))$. We assume that a central limit theorem applies to $\hat{T}(h; k)$, so that in large samples, the t -statistic $[\hat{T}(h; k) - T(f)]/\widehat{\text{se}}(h; k)$ will be approximately normal with variance 1 and mean given by the ratio of bias to standard deviation, $t_f = (E_f[\hat{T}(h; k) - T(f)])/\text{sd}_f(\hat{T}(h; k))$. Since t_f depends on the unknown function f , this ratio is unknown. Note, however, that we can bound $|t_f|$ by the worst-case ratio of bias to standard deviation (bias-sd ratio), $t_{\mathcal{F}} = \sup_{f \in \mathcal{F}} |E_f[\hat{T}(h; k) - T(f)]|/\text{sd}_f(\hat{T}(h; k))$. Therefore, if this bias-sd ratio can be computed up to asymptotically negligible terms, we can construct an honest CI as

$$\hat{T}(h; k) \pm \text{cv}_{1-\alpha}(t) \cdot \widehat{\text{se}}(h; k), \quad (1)$$

where the approximate bias-sd ratio t satisfies $t = t_{\mathcal{F}}(1 + o(1))$, and $\text{cv}_{1-\alpha}(t)$ is the $1 - \alpha$ quantile of the folded normal distribution $|N(t, 1)|$, or, equivalently, the square root of the $1 - \alpha$ quantile of a χ^2 distribution with 1 degree of freedom, and noncentrality parameter t^2 , which is readily available in statistical software. For easy reference, we list these critical values in Table 1 for selected values of t . Because the quantiles of a χ^2 distribution are increasing in its noncentrality parameter, replacing t_f with an upper bound

TABLE 1. Critical values $\text{cv}_{1-\alpha}(\cdot)$.

r	t	α		
		0.01	0.05	0.1
	0.0	2.576	1.960	1.645
6/7	0.408	2.764	2.113	1.777
4/5	0.5	2.842	2.181	1.839
2/3	0.707	3.037	2.362	2.008
1/2	1.0	3.327	2.646	2.284
	1.5	3.826	3.145	2.782
	2.0	4.326	3.645	3.282

Note: Critical values $\text{cv}_{1-\alpha}(t)$ and $\text{cv}_{1-\alpha}(\sqrt{1/r-1})$, for the FLCIs in (1) and (8), corresponding to the $1 - \alpha$ quantiles of the $|N(t, 1)|$ and $|N(\sqrt{1/r-1}, 1)|$ distributions, where t is the bias-sd ratio, and r is the rate exponent. For $t \geq 2$, $\text{cv}_{1-\alpha}(t) \approx t + z_{1-\alpha/2}$ up to 3 decimal places for these values of α .

that is valid for all $f \in \mathcal{F}$ yields a CI that is honest over \mathcal{F} . The CI in (1) is an approximate version of a fixed-length confidence interval (FLCI) studied in [Donoho \(1994\)](#), who replaces $\widehat{\text{se}}(h; k)$ with $\text{sd}_f(\hat{T}(h; k))$ in the definition of this CI, and assumes $\text{sd}_f(\hat{T}(h; k))$ is constant over f , in which case its length will be fixed. We thus refer to CIs of this form as “fixed-length,” even though $\widehat{\text{se}}(h; k)$ is random.

To motivate our main regularity condition (4) below that will facilitate studying the performance of these FLCIs and allow for an easy computation of the bias-sd ratio t , suppose that the standard deviation and the worst-case bias of the estimator $\hat{T}(h; k)$,

$$\overline{\text{bias}}(\hat{T}(h; k)) = \sup_{f \in \mathcal{F}} |E_f \hat{T}(h; k) - T(f)|,$$

scale as powers of h . In particular, suppose that, for some $\gamma_b > 0$, $\gamma_s < 0$, $B(k) > 0$ and $S(k) > 0$,

$$\overline{\text{bias}}(\hat{T}(h; k)) = h^{\gamma_b} MB(k)(1 + o(1)), \quad \text{sd}_f(\hat{T}(h; k)) = h^{\gamma_s} n^{-1/2} S(k)(1 + o(1)), \quad (2)$$

where the $o(1)$ term in the second equality is uniform over $f \in \mathcal{F}$. We show in Appendix B in the Online Supplemental Material that this condition will hold whenever the renormalization heuristics of [Donoho and Low \(1992\)](#) can be formalized. This includes most classical nonparametric problems, such as estimation of a density or a conditional mean, or its derivative, evaluated at a point (which may be a boundary point). In Section 3.2.1, we show that (2) holds with $\gamma_b = p$, and $\gamma_s = -1/2$ under mild regularity conditions when $\hat{T}(h; k)$ is a local polynomial estimator of a conditional mean at a point, and $\mathcal{F}(M)$ consists of functions with p th derivative bounded by M .

REMARK 2.1. The second condition in (2) implies that the standard deviation does not depend on the underlying function f asymptotically. In certain settings, such as density estimation (see Appendix C.1 in the Online Supplemental Material), this may require choosing a localized sequence of parameter spaces \mathcal{F}_n , similar to local asymptotic minimax results in parametric settings (e.g., Section 8.7 in [van der Vaart \(1998\)](#)). While we allow for such dependence, we keep any dependence of \mathcal{F} on n implicit in our notation in the main text. Similarly, the quantities $B(k)$ and $S(k)$ generally depend on \mathcal{F} (which if the parameter space is localized includes the localization point), as well as on other nuisance parameters, such as the variance of the regression errors. To prevent notational clutter, we keep this dependence implicit.

Under (2), we can use the ratio $t = h^{\gamma_b - \gamma_s} MB(k) / (n^{-1/2} S(k))$ of the leading worst-case bias and standard deviation terms to compute the critical value $\text{cv}_{1-\alpha}(t)$ in (1). Analogously to the two-sided case, honest one-sided $1 - \alpha$ CIs based on $\hat{T}(h; k)$ can be constructed by subtracting the standard error times a $1 - \alpha$ quantile of the distribution $\mathcal{N}(t, 1)$. This is asymptotically equivalent to the CI

$$[\hat{T}(h; k) - h^{\gamma_b} MB(k) - z_{1-\alpha} h^{\gamma_s} n^{-1/2} S(k), \infty), \quad (3)$$

which subtracts the maximum bias, in addition to subtracting $z_{1-\alpha}$ times the standard deviation, from $\hat{T}(h; k)$.

REMARK 2.2. One could also form honest two-sided CIs by simply adding and subtracting the worst case bias, in addition to adding and subtracting the standard error times $z_{1-\alpha/2} = \text{cv}_{1-\alpha}(0)$, the $1 - \alpha/2$ quantile of a standard normal distribution, forming the CI as $\hat{T}(h; k) \pm (h^{\gamma_b} MB(k) + z_{1-\alpha/2} \cdot \widehat{\text{se}}(h; k))$. However, since the estimator $\hat{T}(h; k)$ cannot simultaneously have a large positive and a large negative bias, such CI will be conservative, and longer than the CI given in equation (1).

To discuss the optimal choice of bandwidth h and compare efficiency of different kernels k in forming one- and two-sided CIs, and compare the results to the bandwidth and kernel efficiency results for estimation, it will be useful to introduce notation for a generic performance criterion. Let $R(\hat{T})$ denote the worst-case (over \mathcal{F}) performance of \hat{T} according to a given criterion, and let $\tilde{R}(b, s)$ denote the value of this criterion when $\hat{T} - T(f) \sim N(b, s^2)$. For FLCIs, we can take their half-length as the criterion, which leads to

$$R_{\text{FLCI},\alpha}(\hat{T}(h; k)) = \inf\{\chi : P_f(|\hat{T}(h; k) - T(f)| \leq \chi) \geq 1 - \alpha \text{ for all } f \in \mathcal{F}\},$$

$$\tilde{R}_{\text{FLCI},\alpha}(b, s) = \inf\{\chi : P_{Z \sim N(0,1)}(|sZ + b| \leq \chi) \geq 1 - \alpha\} = s \cdot \text{cv}_{1-\alpha}(b/s).$$

To evaluate one-sided CIs, one needs a criterion other than length, which is infinite. A natural criterion is expected excess length, or quantiles of excess length. We focus here on the quantiles of excess length. For CI of the form (3), its worst-case β quantile of excess length is given by $R_{\text{OCI},\alpha,\beta}(\hat{T}(h; k)) = \sup_{f \in \mathcal{F}} q_{f,\beta}(T(f) - \hat{T}(h; k) + h^{\gamma_b} MB(k) + z_{1-\alpha} h^{\gamma_s} n^{-1/2} S(k))$, where $q_{f,\beta}(Z)$ is the β quantile of a random variable Z . The worst-case β quantile of excess length based on an estimator \hat{T} when $\hat{T} - T(f)$ is normal with variance s^2 and bias ranging between $-b$ and b is $\tilde{R}_{\text{OCI},\alpha,\beta}(b, s) = 2b + (z_{1-\alpha} + z_\beta)s$. Finally, to evaluate $\hat{T}(h; k)$ as an estimator we use the maximum root mean squared error (RMSE) under \mathcal{F} as the performance criterion:

$$R_{\text{RMSE}}(\hat{T}) = \sup_{f \in \mathcal{F}} \sqrt{E_f[\hat{T} - T(f)]^2}, \quad \tilde{R}_{\text{RMSE}}(b, s) = \sqrt{b^2 + s^2}.$$

The key regularity condition that we impose on the class of estimators $\hat{T}(h; k)$ is that their performance can be approximated in large samples by the performance of a normally distributed estimator with bias and standard deviation that scale as powers of h ,

$$R(\hat{T}(h; k)) = \tilde{R}(h^{\gamma_b} MB(k), h^{\gamma_s} n^{-1/2} S(k))(1 + o(1)). \tag{4}$$

For the performance criteria above, if the estimator $\hat{T}(h; k)$ satisfies an appropriate central limit theorem, and equation (2) holds, condition (4) will hold so long as the estimator is centered, so that, up to asymptotically negligible terms, its maximum and minimum bias over \mathcal{F} sum to zero, $\sup_{f \in \mathcal{F}} E_f(\hat{T}(h; k) - T(f)) = -\inf_{f \in \mathcal{F}} E_f(\hat{T}(h; k) - T(f))(1 + o(1))$.³ Heuristically, this follows because if $(\hat{T}(h; k) - E_f \hat{T}(h; k))/\text{sd}_f(\hat{T})$ is

³This centering condition holds automatically by a symmetry argument for kernel or local polynomial estimators if f is a conditional mean or a density, $T(f)$ is its value or its derivative at a point, or a regres-

asymptotically $N(0, 1)$, then under (2), $\hat{T}(h; k) - T(f)$ will be in large samples approximately normal, with standard deviation $h^{\gamma_s} n^{-1/2} S(k)$, and mean bounded above and below by $h^{\gamma_b} MB(k)$. In Section 3.2.1, we verify (4) for the problem of estimation of a conditional mean at a point. For estimation of certain smooth nonlinear functionals of the regression function or nonparametric density, including fuzzy regression discontinuity discussed in Section 3, and estimating a bidder valuation in first price auctions discussed in Appendix C.2 of the Online Supplemental Material, moments of the estimator may not exist. In these cases, one can use Theorems B.1 and B.2 in Appendix B in the Online Supplemental Material to verify (4), which only require a weaker version of (2) stated in terms of convergence in distribution rather than moments, so long as one truncates unbounded loss functions.⁴

We also assume that \tilde{R} is homogeneous of degree one,

$$\tilde{R}(tb, ts) = t\tilde{R}(b, s) \quad \text{for all } t > 0. \quad (5)$$

This condition holds for all three criteria considered above. This allows us to simplify the right-hand side of (4). In particular, using the bias-sd ratio $t = h^{\gamma_b - \gamma_s} MB(k) / (n^{-1/2} S(k))$, write the bandwidth as $h = (tn^{-1/2} S(k) / (MB(k)))^{1/(\gamma_b - \gamma_s)}$. Substituting this expression in (4) and using (5) gives

$$\begin{aligned} R(\hat{T}(h; k)) &= \tilde{R}(t^r n^{-r/2} M^{1-r} S(k)^r B(k)^{1-r}, t^{r-1} n^{-r/2} M^{1-r} S(k)^r B(k)^{1-r})(1 + o(1)) \\ &= n^{-r/2} M^{1-r} S(k)^r B(k)^{1-r} t^{r-1} \tilde{R}(t, 1)(1 + o(1)), \end{aligned} \quad (6)$$

where $r = \gamma_b / (\gamma_b - \gamma_s)$. Since the performance criterion converges at the rate $n^{r/2}$ when M is fixed, we refer to r as the rate exponent (this matches the definition in, e.g., [Donoho and Low \(1992\)](#)). We denote the bandwidth choice that minimizes the right-hand side of (6) for a given performance criterion R by $h_R^* = (n^{-1/2} S(k) t_R^* / (MB(k)))^{1/(\gamma_b - \gamma_s)}$, with $t_R^* = \operatorname{argmin}_t t^{r-1} \tilde{R}(t, 1)$, and assume that t_R^* is finite and strictly greater than zero, which is the case for the performance criteria we consider.

The bandwidth choice h_R^* will be asymptotically optimal so long as it is suboptimal to choose a bandwidth sequence h_n such that the bias or the variance dominates asymptotically, which is the case in the settings considered here. For our main results, we assume this directly by assuming that

$$\begin{aligned} M^{r-1} n^{\frac{r}{2}} R(\hat{T}(h_n; k)) &\rightarrow \infty \\ \text{for any } h_n \text{ with } h_n(nM^2)^{\frac{1}{2(\gamma_b - \gamma_s)}} &\rightarrow \infty \text{ or } h_n(nM^2)^{\frac{1}{2(\gamma_b - \gamma_s)}} \rightarrow 0. \end{aligned} \quad (7)$$

sion discontinuity parameter, and \mathcal{F} bounds its derivatives. In other cases, equation (4) will hold when the estimator is recentered by subtracting $\mathfrak{B} = (\sup_{f \in \mathcal{F}} E_f(\hat{T}(h; k) - T(f)) + \inf_{f \in \mathcal{F}} E_f(\hat{T}(h; k) - T(f))) / 2$, or an estimate $\hat{\mathfrak{B}}$ of \mathfrak{B} that is consistent in the sense that $(\hat{\mathfrak{B}} - \mathfrak{B}) / \hat{\sigma}(h; k)$ converges in probability to zero, uniformly over \mathcal{F} . Recentering the estimator in this way improves the estimator's performance under the criteria that we consider.

⁴For evaluating estimators in these cases, we focus on minimizing the limit of the scaled truncated RMSE $\lim_{c \rightarrow \infty} \lim_{n \rightarrow \infty} n^{r/2} M^{r-1} R_{\ell_c}(\hat{T}(h; k))$, where R_{ℓ_c} denotes the worst-case risk under a version of the RMSE that truncates the squared error loss at c^2 . This is equivalent to minimizing the (untruncated) asymptotic RMSE (see Appendix B.1 in the Online Supplemental Material for details). Under this criterion, the RMSE optimal bandwidth defined below and Theorem 2.2 below are not affected by the truncation.

Under this condition, we only need (4) to hold for bandwidth sequences that are of the same order $(nM^2)^{-1/[2(\gamma_b - \gamma_s)]}$ as the optimal bandwidth h_R^* .⁵ Note that optimal bandwidth is of the same order regardless of the performance criterion—the performance criterion only determines the optimal bandwidth constant through t_R^* .

The next theorem collects implications of these derivations for the performance of different kernels. In particular, we consider minimax performance over bandwidth sequences, that is, bandwidth sequences h_n that achieve the asymptotically best possible worst-case performance in large samples in the sense that $M^{r-1}n^{r/2}(R(\hat{T}(h_n; k)) - \inf_{h>0} R(\hat{T}(h; k))) = o(1)$.

THEOREM 2.1. *Let R be a performance criterion with $\tilde{R}(b, s) > 0$ for all $(b, s) \neq 0$. Suppose that equation (4) holds for any bandwidth sequence h_n with*

$$\liminf_{n \rightarrow \infty} h_n (nM^2)^{1/[2(\gamma_b - \gamma_s)]} > 0$$

and

$$\limsup_{n \rightarrow \infty} h_n (nM^2)^{1/[2(\gamma_b - \gamma_s)]} < \infty,$$

and suppose that equations (5) and (7) hold. Define h_R^* and t_R^* as above, and assume that $t_R^* > 0$ is unique and well-defined. Then:

(i) *The asymptotic minimax performance under the kernel k is given by*

$$\begin{aligned} M^{r-1}n^{r/2} \inf_{h>0} R(\hat{T}(h; k)) &= M^{r-1}n^{r/2} R(\hat{T}(h_R^*; k)) + o(1) \\ &= S(k)^r B(k)^{1-r} (t_R^*)^{r-1} \tilde{R}(t_R^*, 1) + o(1). \end{aligned}$$

(ii) *The asymptotic relative efficiency of two kernels k_1 and k_2 is given by*

$$\lim_{n \rightarrow \infty} \frac{\inf_{h>0} R(\hat{T}(h; k_1))}{\inf_{h>0} R(\hat{T}(h; k_2))} = \frac{S(k_1)^r B(k_1)^{1-r}}{S(k_2)^r B(k_2)^{1-r}}.$$

It depends on the rate r but not on the performance criterion R .

(iii) *If we consider two performance criteria R_1 and R_2 satisfying the conditions above, then the limit of the ratio of optimal bandwidths for these criteria is*

$$\lim_{n \rightarrow \infty} \frac{h_{R_1}^*}{h_{R_2}^*} = \left(\frac{t_{R_1}^*}{t_{R_2}^*} \right)^{1/(\gamma_b - \gamma_s)}.$$

⁵In typical settings, a necessary condition for equation (4) to hold is that the optimal bandwidth h_R^* shrinks at a rate such that $(h_R^*)^{-2\gamma_s} n \rightarrow \infty$ and $h_R^* \rightarrow 0$. If M is fixed, this simply requires that $\gamma_b - \gamma_s > 1/2$, which basically amounts to a requirement that $\mathcal{F}(M)$ imposes enough smoothness so that the problem is not degenerate in large samples. If $M = M_n \rightarrow \infty$, then the condition also requires $n^{r/2}M^{r-1} \rightarrow \infty$, so that M does not increase too quickly.

It depends only on γ_b and γ_s and the performance criteria. If (2) holds, the asymptotically optimal bias-sd ratio is given by

$$\lim_{n \rightarrow \infty} \frac{\overline{\text{bias}}(\hat{T}(h_R^*; k))}{\text{sd}_f(\hat{T}(h_R^*; k))} = \underset{t}{\operatorname{argmin}} t^{r-1} \tilde{R}(t, 1) = t_R^*.$$

It depends only on the performance criterion R and rate exponent r .

Part (i) gives the optimal bandwidth formula for a given performance criterion. The performance criterion only determines the optimal bandwidth constant (the optimal bias-sd ratio) t_R^* .

Part (ii) shows that relative kernel efficiency results do not depend on the performance criterion. In particular, known kernel efficiency results under the RMSE criterion such as those in Fan (1993), Cheng, Fan, and Marron (1997), and Fan et al. (1997) apply unchanged to other performance criteria such as length of FLCIs, excess length of one-sided CIs, or expected absolute error.

Part (iii) shows that the optimal bias-sd ratio for a given performance criterion depends on \mathcal{F} only through the rate exponent r , and does not depend on the kernel. The optimal bias-sd ratio for RMSE, FLCI, and OCI, respectively, are

$$t_{\text{RMSE}}^* = \underset{t>0}{\operatorname{argmin}} t^{r-1} \tilde{R}_{\text{RMSE}}(t, 1) = \underset{t>0}{\operatorname{argmin}} t^{r-1} \sqrt{t^2 + 1} = \sqrt{1/r - 1},$$

$$t_{\text{FLCI}}^* = \underset{t>0}{\operatorname{argmin}} t^{r-1} \tilde{R}_{\text{FLCI}, \alpha}(t, 1) = \underset{t>0}{\operatorname{argmin}} t^{r-1} \text{cv}_{1-\alpha}(t), \quad \text{and}$$

$$t_{\text{OCI}}^* = \underset{t>0}{\operatorname{argmin}} t^{r-1} \tilde{R}_{\text{OCI}, \alpha, \beta}(t, 1) = \underset{t>0}{\operatorname{argmin}} t^{r-1} [2t + (z_{1-\alpha} + z_\beta)] = (1/r - 1) \frac{z_{1-\alpha} + z_\beta}{2}.$$

Figures 1 and 2 plot these quantities as a function of r . Note that the optimal bias-sd ratio is larger for FLCIs (at levels $\alpha = 0.05$ and $\alpha = 0.01$) than for RMSE. Since h is increasing in t , it follows that, for FLCI, the optimal bandwidth *oversmooths* relative to the RMSE optimal bandwidth.

REMARK 2.3. Theorem 2.1 does not address whether further efficiency improvements are possible by using estimators that do not fall into the class $\hat{T}(h; k)$, or by using variable length CIs. However, it follows from Donoho (1994) and Armstrong and Kolesár (2018a) that, in typical settings where our results hold, little further improvement is possible. In particular, these papers give efficiency bounds that, applied to our setting, yield asymptotic lower bounds for $R(\hat{T}^*)/R(\hat{T}(h^*; k^*))$, where \hat{T}^* is the optimal estimator or CI among all procedures (for CIs, this includes variable length CIs, with performance measured in terms of expected length), and h^* and k^* are the optimal bandwidth and kernel. These asymptotic lower bounds depend only on the rate exponent r , and so can be used along with the bounds in Theorem 2.1 to obtain the efficiency of a particular kernel and bandwidth relative to the fully optimal procedure.

One can also form FLCIs centered at the estimator that is optimal for different performance criterion R as $\hat{T}(h_R^*; k) \pm \widehat{\text{se}}(h_R^*; k) \cdot \text{cv}_{1-\alpha}(t_R^*)$. The critical value $\text{cv}_{1-\alpha}(t_R^*)$ depends only on the rate exponent r and the performance criterion R . In particular, the

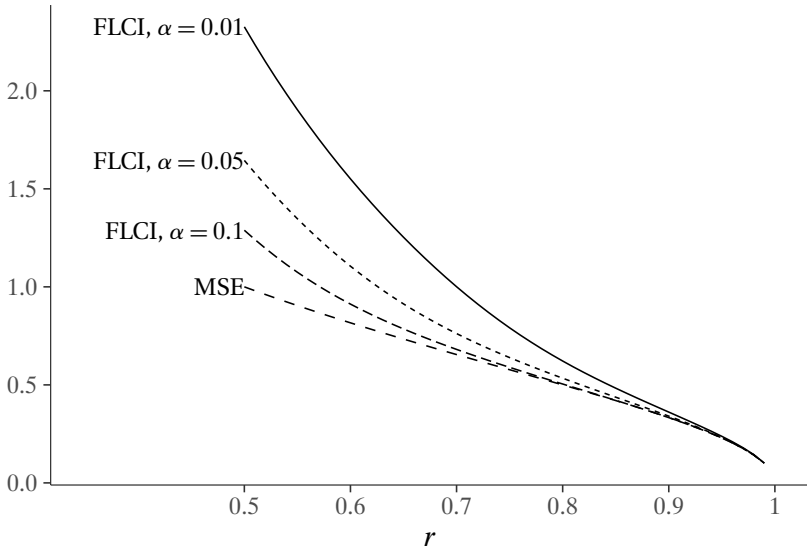


FIGURE 1. Optimal ratio of the worst-case bias to standard deviation for fixed length CIs (FLCI) and maximum MSE (MSE) performance criteria.

CI centered at the RMSE optimal estimator takes this form with $r_{\text{RMSE}}^* = \sqrt{1/r - 1}$, which yields the CI

$$\hat{T}(h_{\text{RMSE}}^*; k) \pm \text{cv}_{1-\alpha}(\sqrt{1/r - 1}) \cdot \hat{\text{se}}(h_{\text{RMSE}}^*; k). \tag{8}$$

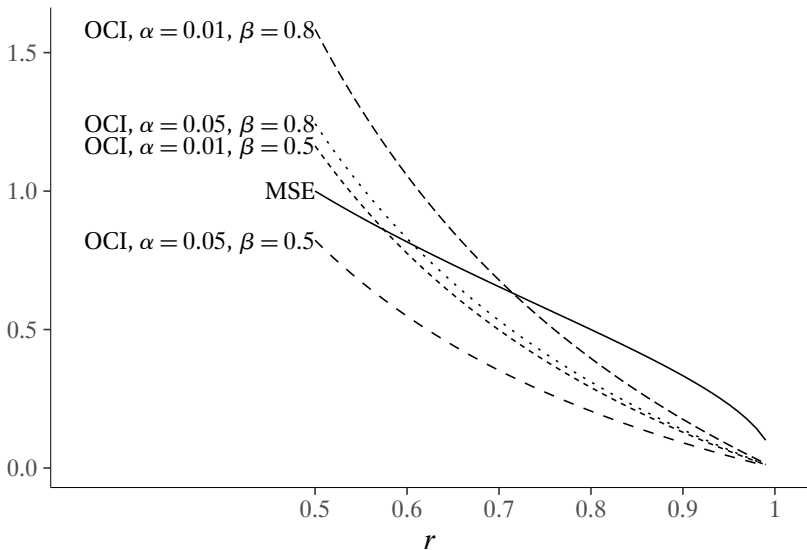


FIGURE 2. Optimal ratio of the worst-case bias to standard deviation for one-sided CIs (OCI), and maximum MSE (MSE) performance criteria.

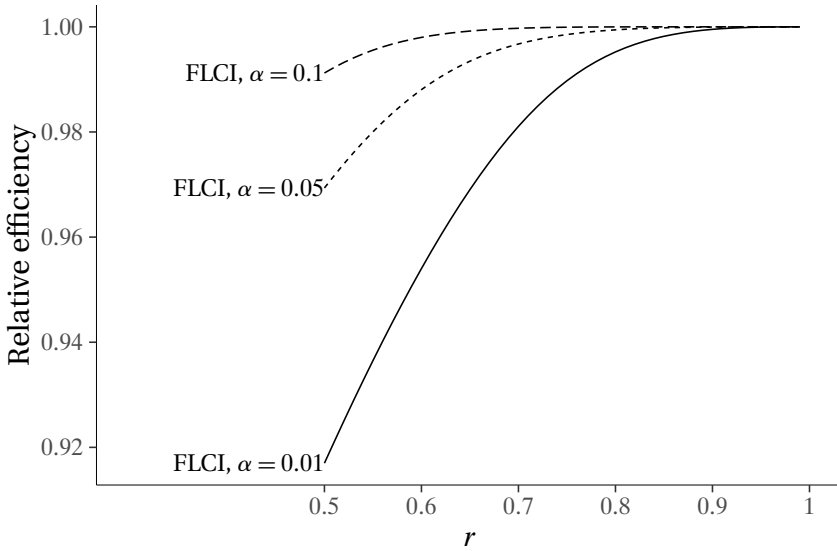


FIGURE 3. Efficiency of fixed-length CIs based on minimax MSE bandwidth relative to fixed-length CIs based on optimal bandwidth.

Table 1 reports this critical value $cv_{1-\alpha}(\sqrt{1/r-1})$ for rate exponents r commonly encountered in practice. By (6), the resulting CI is wider than the one computed using the FLCI optimal bandwidth by a factor of

$$\frac{(t_{\text{FLCI}}^*)^{r-1} \cdot cv_{1-\alpha}(t_{\text{FLCI}}^*)}{(t_{\text{RMSE}}^*)^{r-1} \cdot cv_{1-\alpha}(t_{\text{RMSE}}^*)}. \tag{9}$$

Figure 3 plots this quantity as a function of r . It can be seen from the figure that if $r \geq 4/5$, CIs constructed around the RMSE optimal bandwidth are highly efficient. For example, if $r = 4/5$, to construct an honest 95% FLCI based on an estimator with bandwidth chosen to optimize RMSE, one simply adds and subtracts the standard error multiplied by 2.18 (rather than the usual 1.96 critical value), and the corresponding CI is less than 1% longer than the one with bandwidth chosen to optimize CI length. The next theorem gives a formal statement.

THEOREM 2.2. *Suppose that the conditions of Theorem 2.1 hold for R_{RMSE} and for $R_{\text{FLCI},\tilde{\alpha}}$ for all $\tilde{\alpha}$ in a neighborhood of α . Let $\widehat{se}(h_{\text{RMSE}}^*; k)$ be such that $\widehat{se}(h_{\text{RMSE}}^*; k) / [(h_{\text{RMSE}}^*)^\gamma n^{-1/2} S(k)]$ converges in probability to 1 uniformly over $f \in \mathcal{F}$. Then*

$$\lim_{n \rightarrow \infty} \inf_{f \in \mathcal{F}} P_f(T(f) \in \{\widehat{T}(h_{\text{RMSE}}^*; k) \pm \widehat{se}(h_{\text{RMSE}}^*; k) \cdot cv_{1-\alpha}(\sqrt{1/r-1})\}) = 1 - \alpha.$$

The asymptotic efficiency of this CI relative to the one centered at the FLCI optimal bandwidth, defined as $\lim_{n \rightarrow \infty} \frac{\inf_{h>0} R_{\text{FLCI},\alpha}(\widehat{T}(h;k))}{R_{\text{FLCI},\alpha}(\widehat{T}(h_{\text{RMSE}}^*;k))}$, is given by (9). It depends only on r .

3. APPLICATIONS

In this section, we apply the general results from Section 2 to the problem of inference about a nonparametric regression function at a point, and to regression discontinuity (RD). Readers who are interested only in implementing our CIs in these applications can skip Section 3.2. Appendix C in the Online Supplemental Material discusses two additional applications: estimation of a density at a point, and estimation of a bidder valuation in first-price auctions.

3.1 Setup and estimators

Inference at a point We are interested in inference about a nonparametric regression function f at a point, which we normalize to be zero, so that the parameter of interest is given by $T(f) = f(0)$. We write the nonparametric regression model as

$$y_i = f(x_i) + u_i, \quad i = 1, \dots, n, \quad Eu_i = 0, \quad \text{var}(u_i) = \sigma(x_i). \quad (10)$$

where the design points x_i are nonrandom. We allow the point of interest 0 to lie on the boundary of the support of the design points. We focus on estimating $f(0)$ using a local polynomial estimator of order q with kernel $k(\cdot)$,

$$\hat{T}_q(h; k) = \sum_{i=1}^n w_q^n(x_i; h, k) y_i,$$

where the weights $w_q^n(x_i; h, k)$ are given by

$$w_q^n(x; h, k) = e_1' Q_n^{-1} m_q(x) k(x/h), \quad Q_n = \sum_{i=1}^n k(x_i/h) m_q(x_i) m_q(x_i)'. \quad (11)$$

Here, $m_q(t) = (1, t, \dots, t^q)'$, e_1 is a vector of zeros with 1 in the first position, and h is a bandwidth. Thus, $\hat{T}_q(h; k)$ corresponds to the intercept in a weighted least squares regression of y_i on $(1, x_i, \dots, x_i^q)$ with weights $k(x_i/h)$. Local linear estimators correspond to $q = 1$, and Nadaraya–Watson (local constant) estimators to $q = 0$.

Sharp RD In a sharp RD design, using data from the nonparametric regression model (10), the goal is to estimate the jump in the regression function f at a known cutoff, which we normalize to 0, so that $T(f) = \lim_{x \downarrow 0} f(x) - \lim_{x \uparrow 0} f(x)$. The cutoff determines participation in a binary treatment: units with $x_i \geq 0$ are treated; units with $x_i < 0$ are controls. If the regression functions of potential outcomes are continuous at zero, then $T(f)$ measures the average effect of the treatment for units with $x_i = 0$ (Hahn, Todd, and van der Klaauw (2001)). For brevity, we focus on estimating $T(f)$ based only on local linear regressions: the estimator $\hat{T}(h; k)$ is given by a difference between estimates from two local linear regressions with bandwidth h and kernel k at a boundary point, one for units with nonnegative values running variable x_i , and one for units with negative values of the running variable. The estimator can be written as

$$\hat{T}(h; k) = \sum_{i=1}^n (w_+^n(x; h, k) - w_-^n(x; h, k)) y_i, \quad (12)$$

with the weight w_+^n given by

$$w_+^n(x; h, k) = e_1' Q_{n,+}^{-1} m_1(x) k_+(x/h), \quad k_+(u) = k(u) \mathbf{I}\{u \geq 0\},$$

and $Q_{n,+} = \sum_{i=1}^n k_+(x_i/h) m_1(x_i) m_1(x_i)'$. The weights w_-^n , Gram matrix $Q_{n,-}$ and kernel k_- are defined similarly. Let $\sigma_+^2(x) = \sigma^2(x) \mathbf{I}\{x \geq 0\}$, and $\sigma_-^2(x) = \sigma^2(x) \mathbf{I}\{x < 0\}$.

Fuzzy RD In a fuzzy RD design, the treatment d_i is not entirely determined by whether the running variable x_i exceeds a cutoff. Instead, the cutoff induces a jump in the treatment probability. This fits into our framework if we let $f = (f_1, f_2)$ comprise two regression functions, corresponding to the reduced-form regression of the outcome on the running variable, and the first-stage regression of the treatment on the running variable:

$$\begin{aligned} y_i &= f_1(x_i) + u_{i1}, \\ d_i &= f_2(x_i) + u_{i2}, \end{aligned} \quad i = 1, \dots, n, \quad E u_i = 0, \quad \text{var}(u_i) = \Omega(x_i), \quad (13)$$

with $u_i = (u_{i1}, u_{i2})'$. The parameter of interest is given by the ratio $T(f) = L_1(f)/L_2(f)$ of sharp RD parameters $L_j(f) = \lim_{x \downarrow 0} f_j(x) - \lim_{x \uparrow 0} f_j(x)$ in the reduced-form ($j = 1$) and first-stage regression ($j = 2$). If the regression functions of the potential outcomes and potential treatments are continuous at zero, and a monotonicity condition holds, then $T(f)$ measures the average treatment effect for individuals with $x_i = 0$ who are compliers (see Hahn, Todd, and van der Klaauw (2001)). We consider estimating $T(f)$ by its sample analog, replacing L_1 and L_2 with sharp RD local linear estimates, which are for simplicity assumed to be based on the same bandwidth, $\hat{T}(h; k) = \hat{L}_1(h; k)/\hat{L}_2(h; k)$, where

$$\hat{L}(h; k) = \begin{pmatrix} \hat{L}_1(h; k) \\ \hat{L}_1(h; k) \end{pmatrix} = \sum_i (w_+^n(x; h, k) - w_-^n(x; h, k)) \begin{pmatrix} y_i \\ d_i \end{pmatrix},$$

with the weights w_+^n and w_-^n defined as in (12).

3.2 Theoretical results

We now discuss the conditions under which the key regularity condition (4) holds in each application. We also discuss kernel efficiency results, and gains from imposing global, rather than just local, smoothness on f .

3.2.1 Inference at a point To state the results, it will be convenient to define the equivalent kernel

$$k_q^*(u) = e_1' \left(\int_{\mathcal{X}} m_q(t) m_q(t)' k(t) dt \right)^{-1} m_q(u) k(u), \quad (14)$$

where the integral is over $\mathcal{X} = \mathbb{R}$ if 0 is an interior point, and over $\mathcal{X} = [0, \infty)$ if 0 is a (left) boundary point.

We assume the following conditions on the design points and regression errors u_i .

ASSUMPTION 3.1. For some $d > 0$, the sequence $\{x_i\}_{i=1}^n$ satisfies $\frac{1}{nh_n} \sum_{i=1}^n g(x_i/h_n) \rightarrow d \cdot \int_{\mathcal{X}} g(u) du$ for any bounded function g with finite support and any sequence h_n with $0 < \liminf_n h_n (nM^2)^{1/(2p+1)} < \limsup_n h_n (nM^2)^{1/(2p+1)} < \infty$.

ASSUMPTION 3.2. *The random variables $\{u_i\}_{i=1}^n$ are independent with $Eu_i = 0$, $Eu_i^{2+\eta} \leq 1/\eta$ for some $\eta > 0$, and $\text{var}(u_i) = \sigma^2(x_i)$ for some variance function $\sigma^2(x)$ that is continuous at $x = 0$ with $\sigma^2(0) > 0$.*

Assumption 3.1 requires that the empirical distribution of the design points is smooth around 0. When the support points are treated as random, the constant d typically corresponds to their density at 0.

Because the estimator is linear in y_i , its variance does not depend on f ,

$$\begin{aligned} \text{sd}(\hat{T}_q(h; k))^2 &= \sum_{i=1}^n w_q^n(x_i)^2 \sigma^2(x_i) = \frac{S(k)^2}{nh} (1 + o(1)), \\ S(k) &= \sqrt{\frac{\sigma^2(0) \int_{\mathcal{X}} k_q^*(u)^2 du}{d}}, \end{aligned} \tag{15}$$

where the second equality holds under Assumptions 3.1 and 3.2, as we show in Appendix B.3 in the Online Supplemental Material. The condition on the standard deviation in equation (2) thus holds with $\gamma_s = -1/2$, and $S(k)$ given in the preceding display. Appendix D.3 in the Online Supplemental Material gives the constant $\int_{\mathcal{X}} k_q^*(u)^2 du$ for selected kernels.

On the other hand, the worst-case bias will be driven primarily by the function class \mathcal{F} . We consider inference under two popular function classes. First, the Taylor class of order p ,

$$\mathcal{F}_{T,p}(M) = \left\{ f : \left| f(x) - \sum_{j=0}^{p-1} f^{(j)}(0) x^j / j! \right| \leq M |x|^p / p!, x \in \mathcal{X} \right\}.$$

This class consists of all functions for which the approximation error from a $(p - 1)$ th order Taylor approximation around 0 can be bounded by $\frac{1}{p!} M |x|^p$. It formalizes the idea that the p th derivative of f at zero should be bounded by some constant M . Using this class of functions to derive optimal estimators goes back at least to Legostaeva and Shiryaev (1971), and it underlies much of existing minimax theory concerning local polynomial estimators (see Fan and Gijbels (1996, Chapters 3.4–3.5)).

While analytically convenient, the Taylor class may not be attractive in some empirical settings because it allows f to be nonsmooth and discontinuous away from 0. We therefore also consider inference under Hölder classes (for simplicity, we focus on Hölder classes of integer order)

$$\mathcal{F}_{\text{Hö},p}(M) = \{f : |f^{(p-1)}(x) - f^{(p-1)}(x')| \leq M |x - x'|, x, x' \in \mathcal{X}\}.$$

This class is the closure of the family of p times differentiable functions with the p th derivative bounded by M , uniformly over \mathcal{X} , not just at 0. It formalizes the intuitive notion that f should be p -times differentiable with a bound on the p th derivative. The case $p = 1$ corresponds to the Lipschitz class of functions.

THEOREM 3.1. *Suppose that Assumption 3.1 holds and that $k(\cdot)$ is bounded with bounded support and $q \geq p - 1$. Then, for any bandwidth sequence h_n with $nh_n \rightarrow \infty$ and $0 < \liminf_n h_n(nM^2)^{1/(2p+1)} < \limsup_n h_n(nM^2)^{1/(2p+1)} < \infty$,*

$$\overline{\text{bias}}_{\mathcal{F}_{T,p}(M)}(\hat{T}_q(h_n; k)) = \frac{Mh_n^p}{p!} \mathcal{B}_{p,q}^T(k)(1 + o(1)), \quad \mathcal{B}_{p,q}^T(k) = \int_{\mathcal{X}} |u^p k_q^*(u)| du$$

and

$$\overline{\text{bias}}_{\mathcal{F}_{\text{H\"{o}l},p}(M)}(\hat{T}_q(h_n; k)) = \frac{Mh_n^p}{p!} \mathcal{B}_{p,q}^{\text{H\"{o}l}}(k)(1 + o(1)),$$

$$\mathcal{B}_{p,q}^{\text{H\"{o}l}}(k) = p \int_{t=0}^{\infty} \left| \int_{u \in \mathcal{X}, |u| \geq t} k_q^*(u) (|u| - t)^{p-1} du \right| dt.$$

Thus, the first part of equation (2) holds with $\gamma_b = p$ and $B(k) = \mathcal{B}_{p,q}(k)/p!$, where $\mathcal{B}_{p,q}(k) = \mathcal{B}_{p,q}^{\text{H\"{o}l}}(k)$ for $\mathcal{F}_{\text{H\"{o}l},p}(M)$, and $\mathcal{B}_{p,q}(k) = \mathcal{B}_{p,q}^T(k)$ for $\mathcal{F}_{T,p}(M)$.

If, in addition, Assumption 3.2 holds, then equation (4) holds for the RMSE, FLCI, and OCI performance criteria, with γ_b and $B(k)$ given above and γ_s and $S(k)$ given in equation (15).

The theorem verifies the regularity conditions needed for the results in Section 2, and implies that $r = 2p/(2p + 1)$ for $\mathcal{F}_{T,p}(M)$ and $\mathcal{F}_{\text{H\"{o}l},p}(M)$. If $p = 2$, then we obtain $r = 4/5$. By Theorem 2.1(i), the optimal rate of convergence of a criterion R is $R(\hat{T}(h_R^*; k)) = O((n/M^{1/p})^{-p/(2p+1)})$. As we will see from the relative efficiency calculation below, the optimal order of the local polynomial regression is $q = p - 1$ for the kernels considered here. The theorem allows $q \geq p - 1$, so that we can examine the efficiency of local polynomial regressions that are of order that is too high relative to the smoothness class. Allowing for $q < p - 1$ is not meaningful, as in this case, the maximum bias is infinite.⁶

Under the Taylor class $\mathcal{F}_{T,p}(M)$, the least favorable (bias-maximizing) function is given by $f(x) = M/p! \cdot \text{sign}(w_q^n(x))|x|^p$. In particular, if the weights are not all positive, it will be discontinuous away from the boundary. The first part of Theorem 3.1 then follows by taking the limit of the bias under this function. Assumption 3.1 ensures that this limit is well-defined. Under the Hölder class $\mathcal{F}_{\text{H\"{o}l},p}(M)$, the least favorable function takes the form of a p th order spline. See Appendix B.3 in the Online Supplemental Material for details.

These results imply that given a kernel k and order of a local polynomial q , the RMSE-optimal bandwidth for $\mathcal{F}_{T,p}(M)$ and $\mathcal{F}_{\text{H\"{o}l},p}(M)$ is given by

$$h_{\text{RMSE}}^* = \left(\frac{1}{2pn} \frac{S(k)^2}{M^2 B(k)^2} \right)^{\frac{1}{2p+1}} = \left(\frac{\sigma^2(0) p!^2 \int_{\mathcal{X}} k_q^*(u)^2 du}{2pndM^2 \mathcal{B}_{p,q}(k)^2} \right)^{\frac{1}{2p+1}}, \quad (16)$$

⁶ The smoothness classes $\mathcal{F}_{T,p}(M)$ and $\mathcal{F}_{\text{H\"{o}l},p}(M)$ do not restrict derivatives of order $p - 1$ and lower, so that, in order to achieve a finite worst-case bias, the estimator needs to be unbiased for polynomials of order $p - 1$, which requires $q \geq p - 1$.

where $\mathcal{B}_{p,q}(k) = \mathcal{B}_{p,q}^{\text{Höl}}(k)$ for $\mathcal{F}_{\text{Höl},p}(M)$, and $\mathcal{B}_{p,q}(k) = \mathcal{B}_{p,q}^{\text{T}}(k)$ for $\mathcal{F}_{\text{T},p}(M)$. For kernels given by polynomial functions over their support, k_q^* also has the form of a polynomial, and $\mathcal{B}_{p,q}^{\text{T}}$ and $\mathcal{B}_{p,q}^{\text{Höl}}$ can be computed analytically. Appendix D.3 in the Online Supplemental Material gives these constants for selected kernels.

Kernel efficiency It follows from Theorem 2.1(ii) that the optimal equivalent kernel minimizes $S(k)^r B(k)^{1-r}$, independently of the performance criterion. Under the Taylor class $\mathcal{F}_{\text{T},p}(M)$, this is equivalent to minimizing

$$\left(\int_{\mathcal{X}} k^*(u)^2 du \right)^p \cdot \int_{\mathcal{X}} |u^p k^*(u)| du, \quad (17)$$

The solution to this problem follows from Sacks and Ylvisaker (1978, Theorem 1) (see also Cheng, Fan, and Marron (1997)). We give details of the solution in Appendix D.2 in the Online Supplemental Material. Table 2 compares the asymptotic relative efficiency of local polynomial estimators based on the uniform, triangular, and Epanechnikov kernels to the optimal Sacks–Ylvisaker kernels. Fan et al. (1997) and Cheng, Fan, and Marron (1997) conjectured that minimizing (17) yields a sharp bound on kernel efficiency. It follows from Theorem 2.1(ii) that this conjecture is correct, and Table 2 matches the kernel efficiency bounds in these papers. Table 2 shows that the choice of the kernel does not matter very much, so long as the local polynomial is of the right order. However, if the order is too high, $q > p - 1$, the efficiency can be quite low, even if the bandwidth used was optimal for the function class or the right order, $\mathcal{F}_{\text{T},p}(M)$, especially on the boundary. If the bandwidth picked is optimal for $\mathcal{F}_{\text{T},q-1}(M)$, it will shrink at a lower rate than optimal under $\mathcal{F}_{\text{T},p}(M)$, and the resulting rate of convergence will be lower than r . Consequently, the relative asymptotic efficiency will be zero. A similar point in the context of pointwise asymptotics was made in Sun (2005, Remark 5, p. 8).

The solution to minimizing $S(k)^r B(k)^{1-r}$ under $\mathcal{F}_{\text{Höl},p}(M)$ is only known in special cases. When $p = 1$, the optimal estimator is a local constant estimator based on the triangular kernel. When $p = 2$, the solution is given in Fuller (1961) and Zhao (1997) for

TABLE 2. Relative efficiency of local polynomial estimators for the function class $\mathcal{F}_{\text{T},p}(M)$.

Kernel	Order	Boundary Point			Interior Point		
		$p = 1$	$p = 2$	$p = 3$	$p = 1$	$p = 2$	$p = 3$
Uniform $\mathbb{I}\{ u \leq 1\}$	0	0.9615			0.9615		
	1	0.5724	0.9163		0.9615	0.9712	
	2	0.4121	0.6387	0.8671	0.7400	0.7277	0.9267
Triangular $(1 - u)_+$	0	1			1		
	1	0.6274	0.9728		1	0.9943	
	2	0.4652	0.6981	0.9254	0.8126	0.7814	0.9741
Epanechnikov $\frac{3}{4}(1 - u^2)_+$	0	0.9959			0.9959		
	1	0.6087	0.9593		0.9959	1	
	2	0.4467	0.6813	0.9124	0.7902	0.7686	0.9672

Note: Efficiency is relative to the optimal equivalent kernel k_{SY}^* . The functional $T(f)$ corresponds to the value of f at a point.

TABLE 3. Relative efficiency of local polynomial estimators for the function class $\mathcal{F}_{\text{Hö}l,p}(M)$.

Kernel	Order	Boundary Point			Interior Point		
		$p = 1$	$p = 2$	$p = 3$	$p = 1$	$p = 2$	$p = 3$
Uniform $I\{ u \leq 1\}$	0	0.9615			0.9615		
	1	0.7211	0.9711		0.9615	0.9662	
	2	0.5944	0.8372	0.9775	0.8800	0.9162	0.9790
Triangular $(1 - u)_+$	0	1			1		
	1	0.7600	0.9999		1	0.9892	
	2	0.6336	0.8691	1	0.9263	0.9487	1
Epanechnikov $\frac{3}{4}(1 - u^2)_+$	0	0.9959			0.9959		
	1	0.7471	0.9966		0.9959	0.9949	
	2	0.6186	0.8602	0.9974	0.9116	0.9425	1

Note: For $p = 1, 2$, efficiency is relative to the optimal kernel, for $p = 3$, efficiency is relative to the local quadratic estimator with triangular kernel. The functional $T(f)$ corresponds to the value of f at a point.

the interior point problem, and in Gao (2018) for the boundary point problem. See Appendix D.2 in the Online Supplemental Material for details. When $p \geq 3$, the solution is unknown. Therefore, for $p = 3$, we compute efficiencies relative to a local quadratic estimator with a triangular kernel. Table 3 calculates the resulting efficiencies for local polynomial estimators based on the uniform, triangular, and Epanechnikov kernels. Relative to the class $\mathcal{F}_{T,p}(M)$, the bias constants are smaller: imposing smoothness away from the point of interest helps to reduce the worst-case bias. Furthermore, the loss of efficiency from using a local polynomial estimator of order that is too high is smaller. Finally, local linear regression with a triangular kernel achieves high asymptotic efficiency under both $\mathcal{F}_{T,2}(M)$ and $\mathcal{F}_{\text{Hö}l,2}(M)$, both at the interior and at a boundary, with efficiency at least 97%, giving a theoretical justification to this popular choice in empirical work.

Gains from imposing smoothness globally The Taylor class $\mathcal{F}_{T,p}(M)$, only restricts the p th derivative locally to the point of interest, while the Hölder class $\mathcal{F}_{\text{Hö}l,p}(M)$ restricts the p th derivative globally. How much can one tighten a confidence interval or reduce the RMSE due to this additional smoothness?

It follows from Theorem 3.1 and from arguments underlying Theorem 2.1 that the performance of using a local polynomial estimator of order $p - 1$ with kernel k_H and optimal bandwidth under $\mathcal{F}_{\text{Hö}l,p}(M)$ relative to using a local polynomial estimator of order $p - 1$ with kernel k_T and optimal bandwidth under $\mathcal{F}_{T,p}(M)$ is given by

$$\frac{\inf_{h>0} R_{\mathcal{F}_{\text{Hö}l,p}(M)}(\hat{T}(h; k_H))}{\inf_{h>0} R_{\mathcal{F}_{T,p}(M)}(\hat{T}(h; k_T))} = \left(\frac{\int_{\mathcal{X}} k_{H,p-1}^*(u)^2 du}{\int_{\mathcal{X}} k_{T,p-1}^*(u)^2 du} \right)^{\frac{p}{2p+1}} \left(\frac{\mathcal{B}_{p,p-1}^{\text{Hö}l}(k_H)}{\mathcal{B}_{p,p-1}^T(k_T)} \right)^{\frac{1}{2p+1}} (1 + o(1)), \tag{18}$$

where $R_{\mathcal{F}}(\hat{T})$ denotes the worst-case performance of \hat{T} over \mathcal{F} . If the same kernel is used, the first term equals 1, and the efficiency ratio is determined by the ratio of the

TABLE 4. Gains from imposing global smoothness.

Kernel	Boundary Point			Interior Point		
	$p = 1$	$p = 2$	$p = 3$	$p = 1$	$p = 2$	$p = 3$
Uniform	1	0.855	0.764	1	1	0.848
Triangular	1	0.882	0.797	1	1	0.873
Epanechnikov	1	0.872	0.788	1	1	0.866
Optimal	1	0.906		1	0.995	

Note: The table gives the relative asymptotic risk of local polynomial estimators of order $p - 1$ and a given kernel under the class $\mathcal{F}_{\text{HöL},p}(M)$ relative to the risk under $\mathcal{F}_{T,p}(M)$ given in equation (18). “Optimal” refers to using the optimal kernel under a given smoothness class.

bias constants $\mathcal{B}_{p,p-1}(k)$. Table 4 computes the resulting efficiency gain for common kernels. In general, the gains are greater for larger p , and greater at the boundary. For estimation at a boundary point with $p = 2$, for example, imposing global smoothness of f reduces CI length by about 13–15%, depending on the kernel, and about 10% if the optimal kernel is used.

3.2.2 Sharp regression discontinuity We focus on the most empirically relevant case in which the regression function f is assumed to lie in the class $\mathcal{F}_{\text{HöL},2}(M)$ on either side of the cutoff:

$$f \in \mathcal{F}_{\text{SRD}}(M) = \{f_+(x) \mathbf{I}\{x \geq 0\} - f_-(x) \mathbf{I}\{x < 0\} : f_+, f_- \in \mathcal{F}_{\text{HöL},2}(M)\}.$$

Inference on $T(f)$ is then equivalent to inference on the difference between two regression functions evaluated at boundary points, and the results follow by a slight extension of the results for estimation at a boundary point in Section 3.2.1.

It follows from the results in Section 3.2.1 that if Assumptions 3.1 and 3.2 hold (with the requirement that $\sigma^2(x)$ is continuous 0 replaced by right- and left-continuity of $\sigma_+^2(x)$ and $\sigma_-^2(x)$), then the variance of the estimator does not depend on f and satisfies

$$\begin{aligned} \text{sd}(\hat{T}(h; k))^2 &= \sum_{i=1}^n \tilde{w}^n(x_i)^2 \sigma^2(x_i) = \frac{S(k)^2}{nh} (1 + o(1)), \\ S(k)^2 &= \frac{\int_0^\infty k_1^*(u)^2 du (\sigma_+^2(0) + \sigma_-^2(0))}{d}, \end{aligned}$$

with d defined in Assumption 3.1, and $\tilde{w}^n(x_i) = w_+^n(x_i) + w_-^n(x_i)$. Theorem 3.1 and arguments in Appendix B.3 in the Online Supplemental Material imply that the bias of $\hat{T}(h; k)$ is maximized at $f(x) = -Mx^2/2 \cdot (\mathbf{I}\{x \geq 0\} - \mathbf{I}\{x < 0\})$, so long as the kernel $k(\cdot)$ takes on nonnegative values. The worst-case bias therefore satisfies

$$\overline{\text{bias}}(\hat{T}(h; k)) = -\frac{M}{2} \sum_{i=1}^n \tilde{w}^n(x_i) x_i^2 = Mh^2 B(k) (1 + o(1)), \quad B(k) = -\int_0^\infty u^2 k_1^*(u) du.$$

It follows that for the RMSE, FLCI, and OCI criteria, equation (4) holds with $\gamma_b = 2$, $\gamma_s = -1/2$, and $B(k)$ and $S(k)$ given in the displays above. Thus, the RMSE-optimal bandwidth is given by

$$h_{\text{RMSE}}^* = \left(\frac{\int_0^\infty k_1^*(u)^2 du}{\left(\int_0^\infty u^2 k_1^*(u) du \right)^2} \cdot \frac{\sigma_+^2(0) + \sigma_-^2(0)}{4dnM^2} \right)^{1/5}. \tag{19}$$

The kernel efficiency results are analogous to those in Section 3.2.1.

In principle, one could allow the bandwidths on either side of the cutoff to be different. We show in Appendix D.1 in the Online Supplemental Material, however, that the loss in efficiency resulting from constraining the bandwidths to be the same is quite small unless the ratio of variances on either side of the cutoff, $\sigma_+^2(0)/\sigma_-^2(0)$, is quite large.

3.2.3 Fuzzy regression discontinuity We assume that $f = (f_1, f_2)$ lies in the class $\mathcal{F}_{\text{FRD}}(M_1, M_2) = \mathcal{F}_{\text{SRD}}(M_1) \times \mathcal{F}_{\text{SRD}}(M_2)$, so that both the reduced-form and the first-stage regression functions are assumed to have a bounded second derivative on either side of the cutoff.⁷

Since the estimator is nonlinear, to ensure that (4) holds, it will be necessary to consider a sequence of parameter spaces $\mathcal{F}_{\text{FRD},n}(M_1, M_2)$ localized around a particular value L^* of $L(f) = (L_1(f), L_2(f))'$ with a nonzero jump in the first-stage regression $L_2^* \neq 0$. This allows us to apply a version of the delta method to $\hat{L}(h; k)$. We defer details to Appendix B.4 in the Online Supplemental Material, where we show that under Assumption 3.1 and a version of Assumption 3.2, the distribution of $\hat{T}(h; k) - T(f)$ can in large samples be approximated by a normal distribution with variance

$$\text{avar}(\hat{T}(h; k)) = \frac{S(k)^2}{nh} = \sum_{i=1}^n \frac{s^2(x_i; T(f))}{L_2(f)^2} \tilde{w}^n(x_i; h, k)^2 (1 + o(1)),$$

and mean bounded by

$$\overline{\text{abias}}(\hat{T}(h; k)) = M_1 h^2 B(k) = -\frac{M_1 + |T(f)|M_2}{2|L_2(f)|} \sum_{i=1}^n \tilde{w}^n(x_i; h, k) x_i^2 (1 + o(1)),$$

where $\tilde{w}^n(x_i; h, k) = w_+^n(x_i) + w_-^n(x_i)$, $s^2(x_i; T) = (1, -T)\Omega(x_i)(1, -T)'$,

$$B(k) = -\frac{\int_0^\infty u^2 k_1^*(u) du (1 + |T(f)|M_2/M_1)}{|L_2(f)|},$$

$$S(k)^2 = \frac{\int_0^\infty k_1^*(u)^2 du}{d} \frac{s_+^2(0; T(f)) + s_-^2(0; T(f))}{L_2(f)^2},$$

⁷While we allow the bounds M_1 and M_2 to change with sample size, we assume that their ratio M_1/M_2 is fixed for simplicity.

$s_+^2(0; T) = \lim_{x \downarrow 0} s^2(x; T)$, and $s_-^2(0; T) = \lim_{x \uparrow 0} s^2(x; T)$.

It then follows that for the FLCI, OCI, and a truncated version of the RMSE criterion, equation (4) holds with $M = M_1$, $\gamma_b = 2$, $\gamma_s = -1/2$, and $B(k)$ and $S(k)$ given in the preceding display. The RMSE-optimal bandwidth is therefore given by

$$h_{\text{RMSE}}^* = \left(\frac{\int_0^\infty k_1^*(u)^2 \, du}{\left(\int_0^\infty u^2 k_1^*(u) \, du \right)^2} \cdot \frac{s^2(T(f))}{4dn(M_1 + |T(f)|M_2)} \right)^{1/5}. \quad (20)$$

Since $S(k)$ and $B(k)$ depend on the kernel k through the same quantities as for inference at a boundary point, the kernel efficiency results are analogous to those in Section 3.2.1.

Because the optimal bandwidth depends on $T(f)$, implementing a feasible version of it requires replacing it with an initial estimate. An alternative approach to the construction of two-sided CIs for $T(f)$ that does not require localization or the use of initial estimates is an Anderson and Rubin (1949) style construction studied by Noack and Rothe (2019). In particular, Noack and Rothe (2019) proposed constructing, for each T_0 , an auxiliary CI for the jump in the mean of $y_i - d_i T_0$ at the cutoff, using an approach similar to that we use for inference in sharp RD. The CI for $T(f)$ is then constructed by collecting all T_0 's for which the auxiliary CI contains zero. This approach also has the additional advantage that it can allow for weak identification while it yields asymptotically equivalent CIs under strong identification.⁸ See Noack and Rothe (2019) for a more detailed discussion.

3.3 Practical implementation

We now discuss some practical issues that arise when implementing our CIs for inference at a point, and in sharp and fuzzy RD studied in the previous subsections. To focus the discussion, we consider smoothness classes $\mathcal{F}_{\text{HöL},2}(M)$, $\mathcal{F}_{\text{SRD}}(M)$, and $\mathcal{F}_{\text{FRD}}(M_1, M_2)$ that constrain the second derivative globally, so that, in the discussion below, $p = 2$. In other words, for inference at a point, we assume that the conditional mean f equation (10) is (almost everywhere) twice differentiable with the second derivative bounded by M ; for sharp RD, we assume that that f is twice differentiable on either side of the cutoff, with the second derivative bounded by M ; and for fuzzy RD, we assume that f_1 and f_2 in in equation (13) are twice differentiable on either side of the cutoff, with the second derivative bounded by M_1 and M_2 , respectively. These assumptions imply optimality of the estimators defined in Section 3.1 based on local linear regression ($q = 1$), which is the most popular method in practice; they also imply that both the Epanechnikov and the triangular kernel are nearly optimal.

⁸Because we require that the sequence of parameter spaces $\mathcal{F}_{\text{FRD},n}(M_1, M_2)$ be localized around a value of L^* with $L_2^* \neq 0$, we rule out sequences in which the jump in the first-stage regression is arbitrarily close to zero (the term “weak identification” refers to such sequences). As a result, the CI we propose, unlike the CI proposed by Noack and Rothe (2019), is not honest over the original parameter space $\mathcal{F}_{\text{FRD}}(M_1, M_2)$.

3.3.1 *Choice of M* Appropriate choice of the smoothness constant is key to implementing our method. Since the smoothness classes we consider are convex, the results of [Low \(1997\)](#), [Cai and Low \(2004\)](#) and [Armstrong and Kolesár \(2018a\)](#) imply that, to maintain honesty over the whole function class, a researcher must choose M a priori, rather than attempting to use a data-driven method.⁹ We therefore recommend that, whenever possible, problem-specific knowledge be used to decide what choice of M is reasonable a priori, and that one consider a range of plausible values by way of sensitivity analysis.¹⁰

If one imposes additional restrictions on f that make the parameter space for f non-convex, a data-driven method for choosing M may be feasible.¹¹ In Appendix E in the Online Supplemental Material, we consider a restriction which relates M to a global polynomial approximation to the regression function. In particular, the restriction formalizes the notion that the second derivative in a neighborhood of zero is bounded by the maximum second derivative of a \tilde{p} th order global polynomial approximation. Heuristically, such restriction will hold if the local smoothness of f is no smaller than its smoothness at large scales.

This restriction allows us to calibrate M based on the following rule of thumb. For inference at a point, let $\check{f}(x)$ be an estimate of f based on a global polynomial regression of order \tilde{p} , and let $[x_{\min}, x_{\max}]$ denote the support of x_i . Put $\hat{M}_{\text{ROT}} = \sup_{x \in [x_{\min}, x_{\max}]} |\check{f}^{(\tilde{p})}(x)|$. This rule of thumb is similar to the suggestion of [Fan and Gijbels \(1996, Chapter 4.2\)](#), with the important distinction that their rule of thumb was designed to estimate the pointwise-in- f optimal bandwidth. We discuss the difference between this bandwidth and h_{RMSE}^* in Section 4. In sharp RD, the rule of thumb is analogous, except we define $\check{f}^{(\tilde{p})}(x)$ to be the global polynomial estimate of order \tilde{p} in which the intercept and all coefficients are allowed to be different on either side of the discontinuity (i.e., as regressors, we use $1, x_i, \dots, x_i^{\tilde{p}}$, and their interactions with the indicator $I\{x_i \geq 0\}$). For fuzzy RD, we use an analogous approach to separately calibrate the reduced-form and first-stage smoothness parameters M_1 and M_2 based on the reduced-form and first-stage regressions.

As a default choice, we set $\tilde{p} = p + 2 = 4$. In Appendix E in the Online Supplemental Material, we give a formal analysis of this rule, showing that the resulting CIs are honest and nearly optimal (over a regularity class that imposes the additional restriction f discussed above). In contrast, we expect that calibrating M based on local smoothness estimates may be difficult to justify, since estimating a local derivative of f is a harder problem than the initial problem of estimating its value at a point. We investigate the

⁹These negative results contrast with more positive results for estimation. See, for example, [Lepski \(1990\)](#) who, in the context of estimating the value of the regression function at a point, proposes a data-driven method that automates the choice of both p and M .

¹⁰As is well known, if the final bandwidth choice is influenced by such sensitivity analysis, the resulting CI may undercover, even if the estimator is unbiased. In this case, one can combine our method with the bandwidth snooping adjustment of [Armstrong and Kolesár \(2018b\)](#).

¹¹An alternative to restricting the parameter space is to change the notion of coverage. For example, in the context of constructing confidence bands for a regression function $f(x)$, [Hall and Horowitz \(2013\)](#) proposed bands that have an average coverage property in that the bands achieve coverage of $f(x)$ for a random subset of values of x . This subset may vary with the unknown regression function and the realized sample.

finite-sample performance of FLCIs based on \hat{M}_{ROT} in a Monte Carlo exercise in Section 5.

3.3.2 Computation of RMSE-optimal bandwidth Given a choice of M , one can compute a feasible version \hat{h}_{RMSE}^* of the RMSE-optimal bandwidth by plugging this choice into the expressions (16), (19), and (20), along with consistent estimates of d , and of the variance at 0 (for fuzzy RD, one also needs a preliminary estimate of $T(f)$). In the simulation exercise and empirical application below, we use an alternative approach based on directly minimizing the finite-sample RMSE over the bandwidth h . To describe it, let $\tilde{w}^n(x_i; h, k)$ denote the weights $w_+^n(x_i; h, k)$ given in (11) if the parameter of interest is the conditional mean at a point, and let $\tilde{w}^n(x_i; h, k) = w_+^n(x_i) + w_-^n(x_i)$ if the parameter of interest is the sharp or fuzzy RD parameter.

For inference at a point, or for sharp RD, the finite-sample RMSE takes the form

$$\text{RMSE}(h; M)^2 = \frac{M^2}{4} \left(\sum_{i=1}^n \tilde{w}^n(x_i; h, k) x_i^2 \right)^2 + \sum_{i=1}^n \tilde{w}^n(x_i; h, k) \sigma^2(x_i). \quad (21)$$

Since $\sigma^2(x_i)$ is typically unknown, one needs to replace it by an estimate. For inference at a point, the simplest choice is to use some estimate $\hat{\sigma}^2(x_i) = \hat{\sigma}^2$ that assumes homoskedasticity of the variance function. For sharp RD, one can use the estimate $\hat{\sigma}^2(x_i) = \hat{\sigma}_+^2(0) \mathbf{I}\{x \geq 0\} + \hat{\sigma}_-^2(0) \mathbf{I}\{x < 0\}$, where $\hat{\sigma}_+^2(0)$ and $\hat{\sigma}_-^2(0)$ are some preliminary variance estimates based on observations above and below the cutoff. We use the bandwidth $\hat{h}_{\text{RMSE}, \tilde{M}}^*$ that minimizes equation (21) for $M = \tilde{M}$, the chosen smoothness constant. This method was considered previously in [Armstrong and Kolesár \(2018a\)](#).

Since the estimate in fuzzy RD is nonlinear, its moments, and hence the finite-sample RMSE do not exist. However, one can still employ an analogous approach minimizing the finite-sample analog of the asymptotic RMSE. As the asymptotic bias and the asymptotic standard deviation both scale with the jump in the first-stage regression at the cutoff, $L_2(f)$, this scaling does not affect the optimum, we can equivalently minimize the asymptotic RMSE times $L_2(f)$,

$$\begin{aligned} \text{ARMSE}(h; M_1, M_2)^2 &= \frac{(M_1 + |T(f)|M_2)^2}{4} \left(\sum_{i=1}^n \tilde{w}^n(x_i; h, k) x_i^2 \right)^2 \\ &\quad + \sum_{i=1}^n w_q^n(x_i; h; k)^2 \varsigma^2(x_i; T(f)), \end{aligned}$$

with $\varsigma^2(x; T) = (1, -T)\Omega(x)(1, -T)'$. Since $\Omega(x_i)$ is unknown, one can again replace it with $\hat{\Omega}^2(x_i) = \hat{\Omega}_+^2(0) \mathbf{I}\{x \geq 0\} + \hat{\Omega}_-^2(0) \mathbf{I}\{x < 0\}$, where $\hat{\Omega}_+^2(0)$ and $\hat{\Omega}_-^2(0)$ are some preliminary variance estimates for observations above and below the cutoff. As a preliminary estimate of $T(f)$, one can take the estimate $\hat{T}(\hat{h}_0; k)$, where \hat{h}_0 minimizes the above expression at $T(f) = 0$. One can also use \hat{h}_0 directly as a simple bandwidth selector, which, while not RMSE optimal, has the advantage that it does not depend on the choice of M_2 .

3.3.3 *Construction of FLCIs* Given an estimate \hat{h}_{RMSE}^* of h_{RMSE}^* , such as the estimate $\hat{h}_{\text{RMSE}, \tilde{M}}^*$ discussed above, an honest FLCI can be constructed as

$$\hat{T}(\hat{h}_{\text{RMSE}}^*; k) \pm cv_{1-\alpha}(t) \cdot \widehat{\text{se}}(\hat{h}_{\text{RMSE}}^*; k), \quad (22)$$

where t is an estimate of the bias-sd ratio, and $\widehat{\text{se}}(\hat{h}_{\text{RMSE}}^*; k)$ is an estimate of the standard error. For the standard error, many choices are available in the literature. For inference at a point and sharp RD, the estimator $\hat{T}(\hat{h}_{\text{RMSE}}^*; k)$ is a weighted least squares estimator, and one can directly estimate its finite-sample conditional variance by the nearest neighbor variance estimator considered in [Abadie and Imbens \(2006\)](#) and [Abadie, Imbens, and Zheng \(2014\)](#). Given a bandwidth h , the estimator takes the form

$$\widehat{\text{se}}(h, k)^2 = \sum_{i=1}^n \tilde{w}^n(x_i; h, k)^2 \hat{\sigma}^2(x_i), \quad \hat{\sigma}^2(x_i) = \frac{J}{J+1} \left(y_i - \frac{1}{J} \sum_{j=1}^J y_{j(i)} \right)^2, \quad (23)$$

for some fixed (small) $J \geq 1$, where $j(i)$ denotes the j th closest observation to i (for sharp RD $j(i)$ is only taken among units with the same sign of the running variable.). In contrast, the usual Eicker–Huber–White estimator sets $\hat{\sigma}^2(x_i) = \hat{u}_i^2$, where \hat{u}_i is the regression residual, and it can be shown that this estimator will generally overestimate the conditional variance. For t , one can either use the asymptotic bias-sd ratio $t = 1/2$, or else an estimate of the finite-sample bias-sd ratio $t = -M \sum_{i=1}^n \tilde{w}^n(x_i; \hat{h}_{\text{RMSE}}^*, k) x_i^2 / 2 \widehat{\text{se}}(\hat{h}_{\text{RMSE}}^*, k)$. We use the latter approach in the Monte Carlo and empirical application below. While both approaches are asymptotically equivalent when x_i is continuous, the latter approach has the advantage that it remains valid even when the covariates are discrete.¹²

For fuzzy RD, one can use an analogous approach to estimate the standard error as

$$\widehat{\text{se}}(h, k)^2 = \frac{1}{\hat{L}_2(h; k)^2} \sum_{i=1}^n \tilde{w}^n(x_i; h, k)^2 \hat{\varsigma}^2(x_i, \hat{T}(h; k)),$$

where $\hat{\varsigma}^2(x_i; T) = \frac{J}{J+1} (1, -T) (z_i - \frac{1}{J} \sum_{j=1}^J z_{j(i)}) (z_i - \frac{1}{J} \sum_{j=1}^J z_{j(i)})' (1, -T)'$, $z_i = (y_i, d_i)'$, and $j(i)$ denotes that j th closest observation with the same sign of the running variable. For t , one can use $t = 1/2$, or else the finite-sample analog of the asymptotic bias-sd ratio,¹³ $t = -(\tilde{M}_1 + |\hat{T}| \tilde{M}_2) \cdot \sum_{i=1}^n \tilde{w}^n(x_i; \hat{h}_{\text{RMSE}}^*, k) x_i^2 / 2 \sqrt{\sum_{i=1}^n \hat{\varsigma}^2(x_i; \hat{T}) \tilde{w}^n(x_i; \hat{h}_{\text{RMSE}}^*, k)^2}$.

4. COMPARISON WITH OTHER APPROACHES

In this section, we compare our approach to inference about the parameter $T(f)$ to three other approaches to inference. To make the comparison concrete, we make the comparison in the context of inference about a nonparametric regression function at a point,

¹²See [Armstrong and Kolesár \(2018a\)](#), [Kolesár and Rothe \(2018\)](#), and [Imbens and Wager \(2019\)](#) for a more thorough discussion of the case with discrete covariates.

¹³For inference based on $\hat{T}(\hat{h}_0; k)$, it is necessary to use the finite-sample analog of the bias-sd ratio, since the bandwidth \hat{h}_0 is not RMSE optimal.

discussed in Section 3. The first approach, which we term “conventional,” ignores the potential bias of the estimator and constructs the CI as $\hat{T}_q(h, k) \pm z_{1-\alpha/2}\widehat{\text{se}}(h; k)$. The bandwidth h is typically chosen to minimize the asymptotic mean squared error (MSE) of $\hat{T}_q(h; k)$ under pointwise-in- f (or “pointwise,” for short) asymptotics. We refer to this bandwidth as h_{PT}^* . We discuss the distinction between h_{PT}^* and the bandwidth h_{RMSE}^* in Section 4.1. Under the second approach, undersmoothing, one chooses a sequence of smaller bandwidths, so that in large samples, the bias of the estimator is dominated by its standard error. Finally, in bias correction, one re-centers the conventional CI by subtracting an estimate of the leading bias term from $\hat{T}_q(h; k)$. In Section 4.2, we compare the coverage and length properties of these CIs to the fixed-length CI (FLCI) based on $\hat{T}_q(h_{\text{RMSE}}^*; k)$.

Implementing any of these CIs in practice requires feasible bandwidth and tuning parameter choices. This may require auxiliary assumptions (such as assumptions relating local and global smoothness of f if one picks M using the rule of thumb discussed in Section 3.3.1), which may differ across the methods. For clarity of comparison, we keep implementation issues separate, and focus in this section on a theoretical comparison, assuming any tuning parameters (including the smoothness parameter M) are known. The Monte Carlo exercise in Section 5 below considers their finite-sample performance when the tuning parameters need to be chosen.

4.1 RMSE and pointwise optimal bandwidth

The RMSE optimal bandwidth given in equation (16) seeks to minimize the asymptotic approximation to the maximum RMSE (or, equivalently, MSE) over $f \in \mathcal{F}_{\text{T},p}(M)$ or $f \in \mathcal{F}_{\text{HöL},p}(M)$. In contrast, the bandwidth h_{PT}^* is intended to optimize the MSE at the function f itself. In particular, it minimizes the sum of the leading squared bias and variance terms under pointwise asymptotics for the case $q = p - 1$. It is given by (see, e.g., Fan and Gijbels (1996, equation (3.20)))

$$h_{\text{PT}}^* = \left(\frac{\sigma^2(0)p!^2}{2pndf^{(p)}(0)^2} \frac{\int_{\mathcal{X}} k_q^*(u)^2 du}{\left(\int_{\mathcal{X}} t^p k_q^*(t) dt\right)^2} \right)^{\frac{1}{2p+1}}. \tag{24}$$

Comparing this expression with that for h_{RMSE}^* in equation (16), we see that the pointwise optimal bandwidth replaces M with the p th derivative at zero, $f^{(p)}(0)$, and it replaces $\mathcal{B}_{p,q}(k)$ with $\int_{\mathcal{X}} t^p k_q^*(t) dt$. Note that $\mathcal{B}_{p,q}(k) \geq |\int_{\mathcal{X}} t^p k_q^*(t) dt|$ (this can be seen by noting that the right-hand side corresponds to the bias at the function $f(x) = \pm x^p/p!$, while the left-hand side is the supremum of the bias over functions with p th derivative bounded by (1)). Thus, assuming that $f^{(p)}(0) \leq M$ (this holds by definition for any $f \in \mathcal{F}$ when $\mathcal{F} = \mathcal{F}_{\text{HöL},p}(M)$), we will have $h_{\text{PT}}^*/h_{\text{RMSE}}^* \geq (M/|f^{(p)}(0)|)^{\frac{2}{2p+1}} \geq 1$.

Even though the bandwidth h_{PT}^* is intended to optimize the RMSE at the function f itself, its performance may be arbitrarily bad relative to h_{RMSE}^* at functions for which $f^{(p)}(0)$ is close to zero. For example, consider the function $f(x) = x^{p+1}$ if p is odd, or

$f(x) = x^{p+2}$ if p is even. This is a smooth function with all derivatives bounded on the support of x_i . Since $f^{(p)}(0) = 0$, h_{PT}^* is infinite, and the resulting estimator is a global p th order polynomial least squares estimator. Its RMSE will be poor, since the estimator is not even consistent.¹⁴

To address this problem, plug-in bandwidths that estimate h_{PT}^* include tuning parameters to prevent them from approaching infinity. The RMSE of the resulting estimator at such functions is then determined almost entirely by these tuning parameters. Furthermore, if one uses such a bandwidth as an input to an undersmoothed or bias-corrected CI, the coverage will be determined by these tuning parameters, and can be arbitrarily bad if the tuning parameters allow the bandwidth to be large. Indeed, we find in our Monte Carlo analysis in Section 5 that plug-in estimates of h_{PT}^* used in practice can lead to very poor coverage even when used as a starting point for a bias-corrected or undersmoothed estimator.

4.2 Efficiency and coverage comparison

Let us now consider the efficiency and coverage properties of conventional, undersmoothed, and bias-corrected CIs relative to the FLCI based on $\hat{T}_{p-1}(h_{\text{RMSE}}^*, k)$. To keep the comparison meaningful, and avoid the issues discussed in the previous subsection, we assume these CIs are also based on h_{RMSE}^* , rather than h_{PT}^* (in case of undersmoothing, we assume that the bandwidth is undersmoothed relative to h_{RMSE}^*). Suppose that the smoothness class is either $\mathcal{F}_{\text{T},p}(M)$ or $\mathcal{F}_{\text{Hö},p}(M)$ and denote it by $\mathcal{F}_p(M)$. For concreteness, let $p = 2$, and $q = 1$.

Consider first conventional CIs, given by $\hat{T}_1(h; k) \pm z_{1-\alpha/2} \widehat{\text{se}}(h; k)$. If the bandwidth h equals h_{RMSE}^* , then these CIs are shorter than the 95% FLCIs by a factor of $z_{0.975}/\text{cv}_{0.95}(1/2) = 0.90$. Consequently, their coverage is 92.1% rather than the nominal 95% coverage. At the RMSE-optimal bandwidth, the bias-sd ratio equals $1/2$, so disregarding the bias does not result in severe undercoverage. If one uses a larger bandwidth, however, the bias-sd ratio will be larger, and the undercoverage problem more severe: for example, if the bandwidth is 50% larger than h_{RMSE}^* , so that the bias-sd ratio equals $1/2 \cdot (1.5)^{(5/2)}$, the coverage is only 71.9%.

Second, consider undersmoothing. This amounts to choosing a bandwidth sequence h_n such that $h_n/h_{\text{RMSE}}^* \rightarrow 0$, so that for any fixed M , the bias-sd ratio $t_n = h_n^{\gamma_b - \gamma_s} MB(k)/(n^{-1/2} S(k))$ approaches zero, and the CI $\hat{T}(h^n; k) \pm \text{cv}_{1-\alpha}(0) \widehat{\text{se}}(h_n; k) = \hat{T}(h^n; k) \pm z_{1-\alpha/2} \widehat{\text{se}}(h_n; k)$ will consequently have proper coverage in large samples. However, the CIs shrink at a slower rate than $n^{r/2} = n^{4/5}$, and thus the asymptotic efficiency of the undersmoothed CI relative to the optimal FLCI is zero.

On the other hand, an apparent advantage of the undersmoothed CI is that it appears to avoid specifying the smoothness constant M . However, a more accurate description of undersmoothing is that the bandwidth sequence h_n implicitly chooses a

¹⁴To ensure consistency and finiteness of h_{PT}^* , it is standard to assume that $f^{(p)} \neq 0$. However, the RMSE can still be arbitrarily poor whenever the p th derivative is locally small, but nonzero, and large globally, such as when $f(x) = x^{p+1} + \eta x^p$ for p odd and $f(x) = x^{p+2} + \eta x^p$ if p is even, provided η is sufficiently small.

sequence of smoothness constants $M_n \rightarrow \infty$ such that coverage is controlled under the sequence of parameter spaces $\mathcal{F}_p(M_n)$. We can improve on the coverage and length of the resulting CI by making this sequence explicit and computing an optimal (or near-optimal) FLCI for $\mathcal{F}_p(M_n)$.

To this end, given a sequence h_n , a better approximation to the finite-sample coverage of the CI $\hat{T}(h_n; k) \pm z_{1-\alpha/2} \widehat{\text{se}}(h_n; k)$ over the parameter space $\mathcal{F}_p(M)$ is $P_{Z \sim N(0,1)}(|Z + t_n(M)| \geq z_{1-\alpha/2})$ where $t_n(M) = h_n^{\gamma_b - \gamma_s} MB(k) / (n^{-1/2} S(k))$ is the bias-sd ratio for the given choice of M . This approximation is exact in idealized settings, such as the white noise model discussed in Appendix B in the Online Supplemental Material. For a given level of undercoverage $\eta = \eta_n$, one can then compute M_n as the greatest value of M such that this approximation to the coverage is at least $1 - \alpha - \eta$. In order to trust the undersmoothed CI, one must be convinced of the plausibility of the assumption $f \in \mathcal{F}_p(M_n)$: otherwise, the coverage will be worse than $1 - \alpha - \eta$. This suggests that, in the interest of transparency, one should make this smoothness constant explicit by reporting M_n along with the undersmoothed CI. However, once the sequence M_n is made explicit, a more efficient approach is to simply report an optimal or near-optimal CI for this sequence, either at the coverage level $1 - \alpha - \eta$ (in which case the CI will be strictly smaller than the undersmoothed CI while maintaining the same coverage) or at level $1 - \alpha$ (in which case the CI will have better finite-sample coverage and may also be shorter than the undersmoothed CI).

Finally, let us consider bias correction. It is known that recentering conventional CIs by an estimate of the leading bias term often leads to poor coverage (Hall (1992)). In an important paper, Calonico, Cattaneo, and Titiunik (2014, CCT hereafter) show that the coverage properties of this bias-corrected CI are much better if one adjusts the standard error estimate to account for the variability of the bias estimate, which they call robust bias correction (RBC). For simplicity, consider the case in which the main bandwidth and the pilot bandwidth (used to estimate the bias) are the same, and that the main bandwidth is chosen optimally in that it equals h_{RMSE}^* . In this case, the bias-corrected local linear estimator coincides with a local quadratic estimator. As a result, the RBC procedure in this case amounts to using a local quadratic estimator, but with a bandwidth h_{RMSE}^* , optimal for a local linear estimator. The resulting CI obtains by adding and subtracting $z_{1-\alpha/2}$ times the standard deviation of the estimator.

To ensure that the bias is estimable, the theory of bias correction requires that the conditional mean function is sufficiently smooth, which requires $q < p - 1$ (thus, assuming that f is sufficiently smooth to ensure that the bias of $\hat{T}_1(h; k)$ can be estimated implies that the polynomial order $q = 1$ of the original estimator is not optimal). Suppose, therefore, that the smoothness class is given by $\mathcal{F}_3(M)$ (with $q = 1$, and $h = h_{\text{RMSE}}^*$ still chosen to be MSE optimal for $\mathcal{F}_2(M)$). In this case, the RBC interval can be considered an undersmoothed CI based on a second-order local polynomial estimator. Following the discussion of undersmoothed CIs above, the limiting coverage is $1 - \alpha$ when M is fixed (this matches the pointwise-in- f coverage statements in CCT, which assume the existence of a continuous third derivative in the present context). Due to this undersmoothing, however, the RBC CI shrinks at a slower rate than the optimal CI.

It is also interesting to consider the case when the order $q = 1$ of the local polynomial of the estimator $\hat{T}_1(h_{\text{RMSE}}^*; k)$ is optimal under the maintained smoothness assumption, so that the smoothness class is given by $\mathcal{F}_2(M)$. In this case, the smoothness of the conditional mean function is too low for the bias to be estimable: the bias of the bias-corrected estimator will be of the same order as the bias of the original estimator. Consequently, the estimator will remain asymptotically biased, even after the bias correction. In particular, bias-sd ratio of the estimator is given by

$$t_{\text{RBC}} = (h_{\text{RMSE}}^*)^{5/2} \frac{M \mathcal{B}_{2,2}(k)/2}{\sigma(0) \left(\int_{\mathcal{X}} k_2^*(u)^2 \, du / dn \right)^{1/2}} = \frac{1}{2} \frac{\mathcal{B}_{2,2}(k)}{\mathcal{B}_{2,1}(k)} \left(\frac{\int_{\mathcal{X}} k_1^*(u)^2 \, du}{\int_{\mathcal{X}} k_2^*(u)^2 \, du} \right)^{1/2}. \tag{25}$$

The resulting coverage is given by $\Phi(t_{\text{RBC}} + z_{1-\alpha/2}) - \Phi(t_{\text{RBC}} - z_{1-\alpha/2})$. The RBC interval length relative to the $1 - \alpha$ FLCI around a local linear estimator with the same kernel and minimax MSE bandwidth is the same under both $\mathcal{F}_{T,p}(M)$, and $\mathcal{F}_{\text{Hö}l,p}(M)$, and given by

$$\frac{z_{1-\alpha/2} \left(\int_{\mathcal{X}} k_2^*(u)^2 \, du \right)^{1/2}}{\text{cv}_{1-\alpha}(1/2) \left(\int_{\mathcal{X}} k_1^*(u)^2 \, du \right)^{1/2}} (1 + o(1)). \tag{26}$$

The resulting coverage and relative length is given in Table 5. One can see that although the undercoverage is very mild, (since t_{RBC} is quite low in all cases), the intervals are about 30% longer than the FLCIs around the RMSE bandwidth.

Under the class $\mathcal{F}_{\text{Hö}l,2}(M)$, the RBC intervals are also reasonably robust to using a larger bandwidth: if the bandwidth used is 50% larger than h_{RMSE}^* , so that the bias-sd ratio in equation (25) is larger by a factor of $(1.5)^{5/2}$, the resulting coverage is still at least 93.0% for the kernels considered in Table 5. Under $\mathcal{F}_{T,2}(M)$, using a bandwidth 50%

TABLE 5. Performance of RBC CIs based on h_{RMSE}^* bandwidth for local linear regression under $\mathcal{F}_{T,2}$ and $\mathcal{F}_{\text{Hö}l,2}$.

Kernel	$\mathcal{F}_{T,2}$			$\mathcal{F}_{\text{Hö}l,2}$		
	Length	Coverage	t_{RBC}	Length	Coverage	t_{RBC}
<i>Boundary</i>						
Uniform	1.35	0.931	0.400	1.35	0.948	0.138
Triangular	1.32	0.932	0.391	1.32	0.947	0.150
Epanechnikov	1.33	0.932	0.393	1.33	0.947	0.148
<i>Interior</i>						
Uniform	1.35	0.941	0.279	1.35	0.949	0.086
Triangular	1.27	0.940	0.297	1.27	0.949	0.110
Epanechnikov	1.30	0.940	0.298	1.30	0.949	0.105

Note: Length—CI length relative to 95% FLCI based on a local linear estimator and the same kernel and bandwidth h_{RMSE}^* ; t_{RBC} —ratio of the worst-case bias to standard deviation.

larger than h_{RMSE}^* yields coverage of about 80% on the boundary and 87% in the interior. Thus, depending on the smoothness class, the 95% RBC CI has close to 95% coverage and efficiency loss of about 30%, or exactly 95% coverage at the cost of shrinking at a slower than optimal rate.

Our asymptotic efficiency comparisons focus on minimizing length among CIs with coverage at least $1 - \alpha$ for all $f \in \mathcal{F}$, which follows the usual definition of coverage. One may also consider a criterion that also penalizes CIs that cover “too much,” by placing an upper bound $1 - \underline{\alpha}$ on coverage. For the CIs considered in this paper, the maximum coverage occurs when the bias is zero, and is given by $P_{Z \sim N(0,1)}(|Z| \leq \text{cv}_{1-\alpha}(t)) = 1 - 2\Phi(-\text{cv}_{1-\alpha}(t))$ where t is the asymptotic bias-sd ratio. In particular, when $\mathcal{F} = \mathcal{F}_{\text{T},2}(M)$ or $\mathcal{F} = \mathcal{F}_{\text{HöL},2}(M)$ and the RMSE optimal bandwidth is used, the maximum coverage of a FLCI with 95% (minimum) coverage is $1 - 2\Phi(-2.18) = 0.971$. If one wants the maximum coverage to be smaller, then undersmoothing (or subtracting an estimate of the bias) will be necessary, and Edgeworth expansions may be needed to deal with higher order approximation terms if one wants $\alpha - \underline{\alpha} \rightarrow 0$ quickly enough with the sample size (see Calonico, Cattaneo, and Farrell (2019)). Because, as we discuss above, undersmoothing or bias correction yields longer CIs than the ones we propose, the resulting CIs will be longer than the CIs we propose, which do not penalize “overcoverage.”

5. MONTE CARLO

To study the finite-sample performance of the FLCI that we propose, and compare its performance to other approaches, this section conducts a Monte Carlo analysis of the conditional mean estimation problem considered in Section 3.

We consider Monte Carlo designs with conditional mean functions

$$f_1(x) = \frac{M}{2}(x^2 - 2s(|x| - 0.25)),$$

$$f_2(x) = \frac{M}{2}(x^2 - 2s(|x| - 0.2)^2 + 2s(|x| - 0.5) - 2s(|x| - 0.65)),$$

$$f_3(x) = \frac{M}{2}((x + 1)^2 - 2s(x + 0.2) + 2s(x - 0.2) - 2s(x - 0.4) + 2s(x - 0.7) - 0.92),$$

where $s(x) = (x)_+^2 = \max\{x, 0\}^2$ is the square of the plus function, and $M \in \{2, 6\}$, giving a total of 6 designs. In all cases, x_i is drawn from a uniform distribution with support $[-1, 1]$ (so that the design is random), $u_i \sim N(0, 1/4)$, and the sample size is $n = 500$. Figure 4 plots these designs. The regression function for each design lies in $\mathcal{F}_{\text{HöL},2}(M)$ for the corresponding M . To ensure that our results, discussed below, are not sensitive to the choice of the error distribution or the distribution for the running variable, in Appendix F in the Online Supplemental Material, we also consider designs with x_i drawn from a beta distribution, designs with log-normal and heteroskedastic errors, and designs with different error variance. Finally, we also show in the Appendix that the results remain effectively the same when the function $s(\cdot)$ is replaced by a smooth approximating function.¹⁵

¹⁵The RBC method considered below assumes that the conditional mean function be at least three times continuously differentiable in the neighborhood of 0. Since the functions f_1 , f_2 , and f_3 are not globally

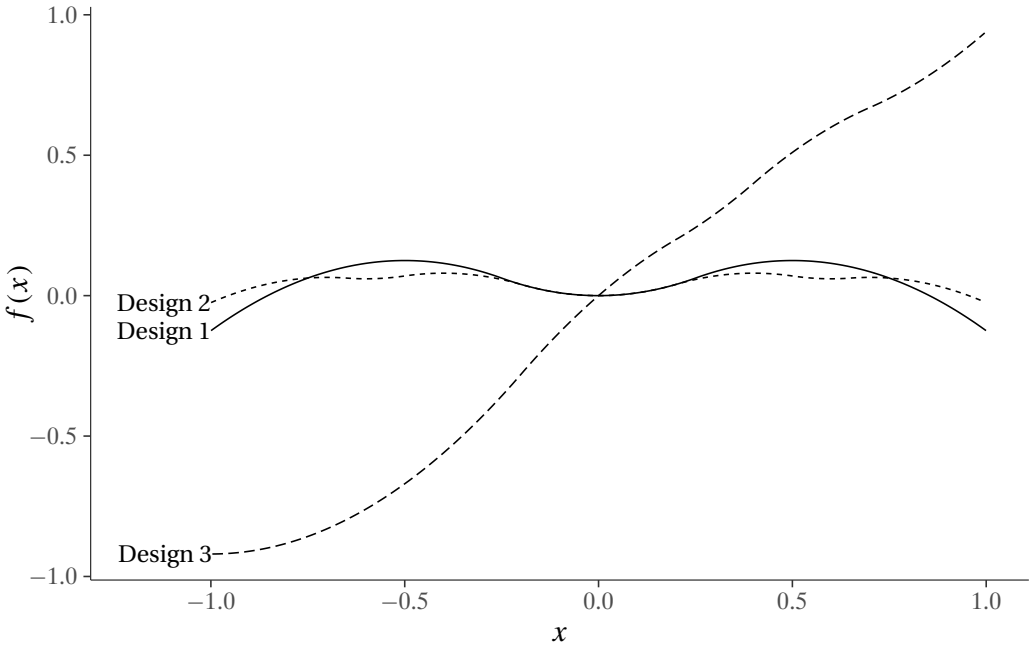


FIGURE 4. Monte Carlo simulation designs 1–3 and $M = 2$.

For each design, we implement the optimal FLCI centered at a local linear estimate with a triangular kernel and MSE optimal bandwidth, as described in Section 3.3, for each choice of $M \in \{2, 6\}$, and with M calibrated using the rule-of-thumb (ROT) described in Section 3.3. The implementations with $M \in \{2, 6\}$ allow us to gauge the effect of using an appropriately calibrated M , compared to a choice of M that is either too conservative or too liberal by a factor of 3. The ROT calibration chooses M automatically, but requires additional conditions in order to have correct coverage (see Section 3.3).

In addition to these FLCIs, we consider seven other CIs (Appendix F in the Online Supplemental Material considers one more method). The first five are different implementations of the robust bias-corrected (RBC) CIs proposed by CCT (discussed in Section 4). Implementing these CIs requires two bandwidth choices: a bandwidth for the local linear estimator, and a pilot bandwidth that is used to construct an estimate of its bias. The first two CIs use bandwidth choices justified by pointwise-in- f asymptotics. The first CI uses a plug-in estimate of h_{pT}^* defined in (24), as implemented by Calonico, Cattaneo, and Farrell (2018), and an analogous estimate for the pilot bandwidth. The second CI, also implemented by Calonico, Cattaneo, and Farrell (2018), uses bandwidth estimates for both bandwidths that optimize the pointwise asymptotic coverage error (CE) among CIs that use usual $z_{1-\alpha/2}$ critical value. This CI can be considered a particular form of undersmoothing. The third CI sets both the pilot bandwidth and the main

three times continuously differentiable, depending on the neighborhood definition, this assumption is arguably violated. The results in the Appendix are nearly identical to those reported here, implying that the performance of the RBC method is not driven by this lack of smoothness.

bandwidth to the plug-in estimate of h_{PT}^* . For the next three CIs, we consider bandwidths justified by uniform-in- f asymptotics. For the fourth and fifth CIs, we set both the main and the pilot bandwidth to h_{RMSE}^* with $M = 2$, and $M = 6$, respectively. For the sixth CI, we set both bandwidths to $\hat{h}_{RMSE, \hat{M}_{ROT}}^*$. Finally, we consider a conventional CI centered at a plug-in bandwidth estimate of h_{PT}^* , using the rule-of-thumb estimator of Fan and Gijbels (1996, Chapter 4.2). All CIs are computed at the nominal 95% coverage level.

Table 6 reports the results. The FLCIs perform well when the correct M is used. As expected, they suffer from undercoverage if M is chosen too small, or suboptimal length when M is chosen too large. The ROT choice of M appears to do a reasonable job of having good coverage and length in these designs without requiring knowledge of the true smoothness constant. However, as discussed in Section 3.3, this ROT choice imposes additional restrictions on the parameter space, so one must take care in extrapolating these results to other designs.

As predicted by the theory in Section 4, the RBC CIs also have good coverage when implemented using the h_{RMSE}^* bandwidth, and they are less sensitive to the choice of M than the corresponding FLCIs, at the expense of being on average about 25% longer. RBC CIs with bandwidth given by $\hat{h}_{RMSE, \hat{M}_{ROT}}^*$ also achieve good coverage, but they are again about 25% longer than the corresponding FLCIs.

The CIs based on bandwidths justified by pointwise-in- f asymptotics (rows 1, 2, 3, and 7 for each design in the table) all have very poor coverage for at least one of the designs. Our analysis in Section 4 suggests that this is due to the tuning parameter choices required by these bandwidths. Indeed, looking at the average of the bandwidth over the Monte Carlo draws (also reported in Table 6), it can be seen that the bandwidths tend to be much larger than those that estimate h_{RMSE}^* . This is even the case for the CE bandwidth, which is intended to minimize coverage errors.

Overall, the Monte Carlo analysis suggests that default approaches to nonparametric CI construction (bias-correction or undersmoothing relative to plug-in bandwidths) can lead to severe undercoverage when implemented using bandwidths justified by pointwise-in- f asymptotics. Bias-corrected CIs such as the one proposed by CCT can have good coverage if one starts from the minimax RMSE bandwidth, although they will be wider than FLCIs proposed in this paper.

6. EMPIRICAL ILLUSTRATION

To illustrate the implementation of feasible versions of the CIs (22), we use a subset of the dataset from Ludwig and Miller (2007).

In 1965, when the Head Start federal program launched, the Office of Economic Opportunity provided technical assistance to the 300 poorest counties in the United States to develop Head Start funding proposals. Ludwig and Miller (2007) used this cutoff in technical assistance to look at intent-to-treat effects of the Head Start program on a variety of outcomes using as a running variable the county's poverty rate relative to the poverty rate of the 300th poorest county (which had poverty rate equal to approximately 59.2%). We focus here on their main finding, the effect on child mortality due to causes addressed as part of Head Start's health services. See Ludwig and Miller (2007) for a

TABLE 6. Monte Carlo simulation: inference at a point.

Method	Bandwidth	$M = 2$					$M = 6$				
		Bias	SE	$E[h]$	Cov	RL	Bias	SE	$E_m[h]$	Cov	RL
<i>Design 1</i>											
RBC	$h = \hat{h}_{PT}^*, b = \hat{b}_{PT}^*$	0.063	0.035	0.75	55.6	0.73	0.157	0.036	0.62	0.1	0.61
RBC	$h = \hat{h}_{CE}, b = \hat{b}_{CE}$	0.030	0.041	0.45	85.8	0.85	0.059	0.045	0.34	72.4	0.76
RBC	$h = b = \hat{h}_{PT}^*$	0.025	0.042	0.75	93.1	0.88	0.042	0.047	0.62	89.1	0.78
RBC	$h = b = \hat{h}_{RMSE,2}^*$	0.001	0.061	0.36	94.5	1.27	0.002	0.061	0.36	94.5	1.01
RBC	$h = b = \hat{h}_{RMSE,6}^*$	0.000	0.076	0.23	94.2	1.58	0.000	0.075	0.23	94.2	1.26
RBC	$h = b = \hat{h}_{RMSE, \hat{M}_{ROT}}^*$	0.000	0.078	0.22	93.9	1.64	0.000	0.097	0.14	93.4	1.63
Conventional	$\hat{h}_{PT,ROT}^*$	0.032	0.036	0.56	76.6	0.76	0.049	0.046	0.31	77.4	0.77
FLCI, $M = 2$	$\hat{h}_{RMSE,2}^*$	0.021	0.043	0.36	94.9	1.00	0.065	0.043	0.36	75.2	0.80
FLCI, $M = 6$	$\hat{h}_{RMSE,6}^*$	0.009	0.054	0.23	96.6	1.25	0.028	0.053	0.23	94.7	1.00
FLCI, $M = \hat{M}_{ROT}$	$\hat{h}_{RMSE, \hat{M}_{ROT}}^*$	0.008	0.056	0.22	95.6	1.29	0.010	0.069	0.14	96.3	1.30
<i>Design 2</i>											
RBC	$h = \hat{h}_{PT}^*, b = \hat{b}_{PT}^*$	0.043	0.035	0.77	75.9	0.72	0.129	0.035	0.77	4.6	0.58
RBC	$h = \hat{h}_{CE}, b = \hat{b}_{CE}$	0.028	0.040	0.49	87.4	0.83	0.074	0.041	0.44	54.1	0.69
RBC	$h = b = \hat{h}_{PT}^*$	0.026	0.041	0.77	90.9	0.87	0.077	0.042	0.77	53.0	0.70
RBC	$h = b = \hat{h}_{RMSE,2}^*$	0.002	0.061	0.36	94.5	1.27	0.006	0.061	0.36	94.4	1.01
RBC	$h = b = \hat{h}_{RMSE,6}^*$	0.000	0.076	0.23	94.2	1.58	0.000	0.075	0.23	94.2	1.26
RBC	$h = b = \hat{h}_{RMSE, \hat{M}_{ROT}}^*$	0.001	0.068	0.30	94.0	1.43	0.000	0.083	0.20	93.8	1.38
Conventional	$\hat{h}_{PT,ROT}^*$	0.032	0.032	0.78	74.4	0.67	0.073	0.040	0.44	53.0	0.66
FLCI, $M = 2$	$\hat{h}_{RMSE,2}^*$	0.020	0.043	0.36	95.1	1.00	0.061	0.043	0.36	78.1	0.80
FLCI, $M = 6$	$\hat{h}_{RMSE,6}^*$	0.009	0.054	0.23	96.6	1.25	0.028	0.053	0.23	94.7	1.00
FLCI, $M = \hat{M}_{ROT}$	$\hat{h}_{RMSE, \hat{M}_{ROT}}^*$	0.013	0.048	0.30	94.3	1.13	0.020	0.059	0.20	94.3	1.10
<i>Design 3</i>											
RBC	$h = \hat{h}_{PT}^*, b = \hat{b}_{PT}^*$	-0.043	0.035	0.77	75.7	0.72	-0.123	0.035	0.74	9.9	0.59
RBC	$h = \hat{h}_{CE}, b = \hat{b}_{CE}$	-0.026	0.040	0.49	88.1	0.83	-0.063	0.043	0.43	64.2	0.71
RBC	$h = b = \hat{h}_{PT}^*$	-0.024	0.042	0.77	90.8	0.87	-0.066	0.043	0.74	60.3	0.71
RBC	$h = b = \hat{h}_{RMSE,2}^*$	-0.002	0.061	0.36	94.5	1.27	-0.007	0.061	0.36	94.4	1.01
RBC	$h = b = \hat{h}_{RMSE,6}^*$	0.000	0.076	0.23	94.2	1.58	0.000	0.075	0.23	94.2	1.26
RBC	$h = b = \hat{h}_{RMSE, \hat{M}_{ROT}}^*$	0.000	0.074	0.25	94.2	1.54	0.000	0.092	0.16	93.6	1.54
Conventional	$\hat{h}_{PT,ROT}^*$	-0.032	0.033	0.72	74.7	0.69	-0.065	0.042	0.39	62.0	0.70
FLCI, $M = 2$	$\hat{h}_{RMSE,2}^*$	-0.020	0.043	0.36	95.0	1.00	-0.060	0.043	0.36	78.1	0.80
FLCI, $M = 6$	$\hat{h}_{RMSE,6}^*$	-0.009	0.054	0.23	96.5	1.25	-0.027	0.053	0.23	94.7	1.00
FLCI, $M = \hat{M}_{ROT}$	$\hat{h}_{RMSE, \hat{M}_{ROT}}^*$	-0.010	0.052	0.25	95.6	1.22	-0.013	0.065	0.16	96.1	1.22

Note: Legend: SE—average standard error; $E[h]$ —average (over Monte Carlo draws) bandwidth; Cov—coverage of CIs (in %); RL—relative (to optimal FLCI) length. *Bandwidth descriptions:* \hat{h}_{PT}^* —plugin estimate of pointwise MSE optimal bandwidth (bw); \hat{b}_{PT}^* —analog for estimate of the bias; \hat{h}_{CE} —plugin estimate of coverage error optimal bw; \hat{b}_{CE} —analog for estimate of the bias; The implementation of Calonico, Cattaneo, and Farrell (2018) is used for all four bws. $\hat{h}_{RMSE,2}^*$, $\hat{h}_{RMSE,6}^*$ —RMSE optimal bw, assuming $M = 2$, and $M = 6$, respectively. $\hat{h}_{PT,ROT}^*$ —Fan and Gijbels (1996) rule of thumb; $\hat{h}_{RMSE, \hat{M}_{ROT}}^*$ —RMSE optimal bw, using rule-of-thumb for M . 50,000 Monte Carlo draws.

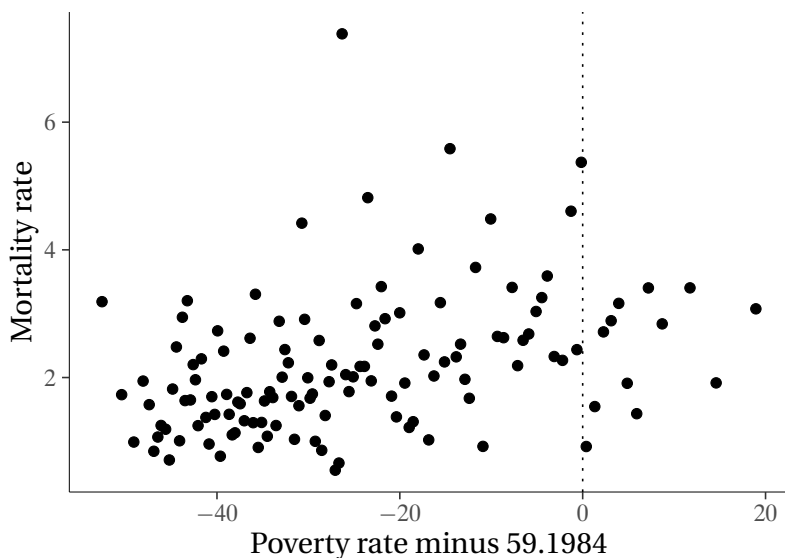


FIGURE 5. Average county mortality rate per 100,000 for children aged 5–9 over 1973–83 due to causes addressed as part of Head Start’s health services (labeled “Mortality rate”) plotted against poverty rate in 1960 relative to the 300th poorest county. Each point corresponds to an average for 25 counties. Data are from Ludwig and Miller (2007).

detailed description of this variable. Relative to the dataset used in Ludwig and Miller (2007), we remove one duplicate entry and one outlier, which after discarding counties with partially missing data leaves us with 3103 observations, with 294 of them above the poverty cutoff.

Figure 5 plots the data (to reduce the noise in the outcome variable, we plot bin averages of size 25). To estimate the discontinuity in mortality rates, Ludwig and Miller (2007) used a uniform kernel¹⁶ and consider bandwidths equal to 9, 18, and 36. This yields point estimates equal to -1.90 , -1.20 , and -1.11 , respectively, which are large effects given that the average mortality rate for counties not receiving technical assistance was 2.15 per 100,000. The p -values reported in the paper, based on bootstrapping the t -statistic (which ignores any potential bias in the estimates), are 0.036, 0.081, and 0.027. The standard errors for these estimates, obtained using the nearest neighbor method (with $J = 3$) are 1.04, 0.70, and 0.52.

These bandwidth choices are optimal in the sense that they minimize the RMSE expression (21) if $M = 0.040$, 0.0074, and 0.0014, respectively. Thus, for these bandwidths to be optimal, one has to be very optimistic about the smoothness of the regression function. In comparison, the rule of thumb method for estimating M discussed in Section 3.3 yields $\hat{M}_{\text{ROT}} = 0.299$, implying h_{RMSE}^* estimate 4.0, and the point estimate -3.17 . For these smoothness parameters, the critical values based on the finite-sample bias-sd ratio are

¹⁶Ludwig and Miller (2007) stated that the estimates were obtained using a triangular kernel. However, due to a bug in the code, the results reported in the paper were actually obtained using a uniform kernel.

given by 2.165, 2.187, 2.107, and 2.202, respectively, which is very close to the asymptotic value $cv_{0.95}(1/2) = 2.181$. The resulting 95% confidence intervals are given by

$$(-4.143, 0.353), \quad (-2.720, 0.323), \quad (-2.215, -0.013), \quad \text{and} \quad (-6.352, 0.010),$$

respectively. The p -values based on these estimates are given by 0.100, 0.125, 0.047, and 0.051. These p -values are larger than those reported in the paper, as they take into account the potential bias of the estimates.

Using a triangular kernel helps to tighten the confidence intervals by a few percentage points in length, as predicted by the relative asymptotic efficiency results from Table 3, yielding

$$(-4.138, 0.187), \quad (-2.927, 0.052), \quad (-2.268, -0.095), \quad \text{and} \quad (-5.980, -0.322).$$

The underlying optimal bandwidths are given by 11.6, 23.1, 45.8, and 4.9, respectively. The p -values associated with these estimates are 0.074, 0.059, 0.033, and 0.028, tightening the p -values based on the uniform kernel.

These results indicate that unless one is very optimistic about the smoothness of the regression function, the uncertainty associated with the magnitude of the effect of Head Start assistance on child mortality is much higher than originally reported. This is due mainly to the relatively large bandwidths used by [Ludwig and Miller \(2007\)](#), which imply an optimistic bound on the smoothness of the regression function if we assume that such bandwidths are close to optimal for MSE. Interestingly, while the more conservative smoothness bound in our benchmark specification leads to much wider CIs, the point estimate is larger in magnitude, so that one still finds a statistically significant effect at a 5 or 10% level, depending on the kernel.

APPENDIX A: PROOFS OF THEOREMS IN SECTION 2

A.1 Proof of Theorem 2.1

Parts (ii) and (iii) follow from part (i) and simple calculations. To prove part (i), note that, if it did not hold, there would be a bandwidth sequence h_n such that

$$\liminf_{n \rightarrow \infty} M^{r-1} n^{r/2} R(\hat{T}(h_n; k)) < S(k)^r B(k)^{1-r} \inf_t t^{r-1} \tilde{R}(t, 1).$$

By equation (7), the bandwidth sequence h_n must satisfy $\liminf_{n \rightarrow \infty} h_n (nM^2)^{1/[2(\gamma_b - \gamma_s)]} > 0$ and $\limsup_{n \rightarrow \infty} h_n (nM^2)^{1/[2(\gamma_b - \gamma_s)]} < \infty$. Thus, by equation (6),

$$M^{r-1} n^{r/2} R(\hat{T}(h_n; k)) = S(k)^r B(k)^{1-r} t_n^{r-1} \tilde{R}(t_n, 1) + o(1),$$

where $t_n = h_n^{\gamma_b - \gamma_s} B(k) / (n^{-1/2} S(k))$. This contradicts the display above.

A.2 Proof of Theorem 2.2

The second statement (relative efficiency) is immediate from (6). For the first statement (coverage), fix $\varepsilon > 0$ and let $sd_n = n^{-1/2} (h_{\text{RMSE}}^*)^{\gamma_s} S(k)$ so that $sd_n / \widehat{\text{se}}(h_{\text{RMSE}}^*; k) \xrightarrow{P} 1$ uni-

formly over $f \in \mathcal{F}$. Note that, by Theorem 2.1 and the fact that $t_{\text{RMSE}}^* = \sqrt{1/r - 1}$,

$$\tilde{R}_{\text{FLCI}, \alpha + \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k)) = \text{sd}_n \cdot \text{cv}_{1-\alpha-\varepsilon}(\sqrt{1/r - 1})(1 + o(1)),$$

and similarly for $\tilde{R}_{\text{FLCI}, \alpha - \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k))$. Since $\text{cv}_{1-\alpha}(\sqrt{1/r - 1})$ is strictly decreasing in α , it follows that there exists $\eta > 0$ such that, with probability approaching 1 uniformly over $f \in \mathcal{F}$,

$$\begin{aligned} R_{\text{FLCI}, \alpha + \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k)) &< \widehat{\text{se}}(\hat{T}(h_{\text{RMSE}}^*; k)) \cdot \text{cv}_{1-\alpha}(\sqrt{1/r - 1}) \\ &< (1 - \eta)R_{\text{FLCI}, \alpha - \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k)). \end{aligned}$$

Thus,

$$\begin{aligned} \liminf_n \inf_{f \in \mathcal{F}} P(T(f) \in \{\hat{T}(h_{\text{RMSE}}^*; k) \pm \widehat{\text{se}}(\hat{T}(h_{\text{RMSE}}^*; k)) \cdot \text{cv}_{1-\alpha}(\sqrt{1/r - 1})\}) \\ \geq \liminf_n \inf_{f \in \mathcal{F}} P(T(f) \in \{\hat{T}(h_{\text{RMSE}}^*; k) \pm R_{\text{FLCI}, \alpha + \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k))\}) \geq 1 - \alpha - \varepsilon, \end{aligned}$$

and

$$\begin{aligned} \limsup_n \inf_{f \in \mathcal{F}} P(T(f) \in \{\hat{T}(h_{\text{RMSE}}^*; k) \pm \widehat{\text{se}}(\hat{T}(h_{\text{RMSE}}^*; k)) \cdot \text{cv}_{1-\alpha}(\sqrt{1/r - 1})\}) \\ \leq \limsup_n \inf_{f \in \mathcal{F}} P(T(f) \in \{\hat{T}(h_{\text{RMSE}}^*; k) \pm R_{\text{FLCI}, \alpha - \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k))(1 - \eta)\}) \leq 1 - \alpha + \varepsilon, \end{aligned}$$

where the last inequality follows by definition of $R_{\text{FLCI}, \alpha - \varepsilon}(\hat{T}(h_{\text{RMSE}}^*; k))$. Taking $\varepsilon \rightarrow 0$ gives the result.

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