



FICO™ Xpress  
Optimization Suite

Xpress Knitro Controls Reference  
Manual

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Xpress-Knitro

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## CHAPTER 1

# Xpress Knitro Control Parameters

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This manual provides a full list of the controls accepted by Xpress for setting Knitro parameters. KNITRO has a great number and variety of user option settings and although it tries to choose the best settings by default, often significant performance improvements can be realized by choosing some non-default option settings.

XKTR_PARAM_ALGORITHM	Indicates which algorithm to use to solve the problem	p. 8
XKTR_PARAM_BAR_DIRECTINTERVAL	Controls the maximum number of consecutive conjugate gradient (CG) steps before KNITRO will try to enforce that a step is taken using direct linear algebra.	p. 8
XKTR_PARAM_BAR_FEASIBLE	Specifies whether special emphasis is placed on getting and staying feasible in the interior-point algorithms.	p. 8
XKTR_PARAM_BAR_FEASMODETOL	Specifies the tolerance in equation that determines whether KNITRO will force subsequent iterates to remain feasible.	p. 4
XKTR_PARAM_BAR_INITMU	Specifies the initial value for the barrier parameter : used with the barrier algorithms. This option has no effect on the Active Set algorithm.	p. 4
XKTR_PARAM_BAR_INITPT	Indicates whether an initial point strategy is used with barrier algorithms.	p. 9
XKTR_PARAM_BAR_MAXBACKTRACK	Indicates the maximum allowable number of backtracks during the linesearch of the Interior/Direct algorithm before reverting to a CG step.	p. 9
XKTR_PARAM_BAR_MAXCROSSIT	Specifies the maximum number of crossover iterations before termination.	p. 9
XKTR_PARAM_BAR_MAXREFACTOR	Indicates the maximum number of refactorizations of the KKT system per iteration of the Interior/Direct algorithm before reverting to a CG step.	p. 10
XKTR_PARAM_BAR_MURULE	Indicates which strategy to use for modifying the barrier parameter mu in the barrier algorithms.	p. 10
XKTR_PARAM_BAR_PENCONS	Indicates whether a penalty approach is applied to the constraints.	p. 11
XKTR_PARAM_BAR_PENRULE	Indicates which penalty parameter strategy to use for determining whether or not to accept a trial iterate.	p. 11
XKTR_PARAM_BAR_SWITCHRULE	Indicates whether or not the barrier algorithms will allow switching from an optimality phase to a pure feasibility phase.	p. 12

XKTR_PARAM_DELTA	Specifies the initial trust region radius scaling factor used to determine the initial trust region size.	p. 4
XKTR_PARAM_DELTA	Specifies the initial trust region radius scaling factor used to determine the initial trust region size.	p. 4
XKTR_PARAM_FEASTOL	Specifies the final relative stopping tolerance for the feasibility error. p. 5	
XKTR_PARAM_FEASTOLABS	Specifies the final absolute stopping tolerance for the feasibility error.	p. 5
XKTR_PARAM_GRADOPT	Specifies how to compute the gradients of the objective and constraint functions.	p. 12
XKTR_PARAM_HESSOPT	Specifies how to compute the (approximate) Hessian of the Lagrangian.	p. 12
XKTR_PARAM_HONORBNDs	Indicates whether or not to enforce satisfaction of simple variable bounds throughout the optimization.	p. 13
XKTR_PARAM_INFEASTOL	Specifies the (relative) tolerance used for declaring infeasibility of a model.	p. 5
XKTR_PARAM_LMSIZE	Specifies the number of limited memory pairs stored when approximating the Hessian using the limited-memory quasi-Newton BFGS option.	p. 13
XKTR_PARAM_MAXCGIT	Specifies the number of limited memory pairs stored when approximating the Hessian using the limited-memory quasi-Newton BFGS option.	p. 14
XKTR_PARAM_MAXIT	Specifies the maximum number of iterations before termination.	p. 14
XKTR_PARAM_MIP_BRANCHRULE	Specifies which branching rule to use for MIP branch and bound procedure.	p. 14
XKTR_PARAM_MIP_GUB_BRANCH	Specifies whether or not to branch on generalized upper bounds (GUBs).	p. 14
XKTR_PARAM_MIP_HEURISTIC	Specifies which MIP heuristic search approach to apply to try to find an initial integer feasible point.	p. 15
XKTR_PARAM_MIP_HEURISTIC_MAXIT	Specifies the maximum number of iterations to allow for MIP heuristic, if one is enabled.	p. 15
XKTR_PARAM_MIP_IMPLICATNS	Specifies whether or not to add constraints to the MIP derived from logical implications.	p. 15
XKTR_PARAM_MIP_INTEGERTOL	This value specifies the threshold for deciding whether or not a variable is determined to be an integer.	p. 5
XKTR_PARAM_MIP_INTGAPABS	The absolute integrality gap stop tolerance for MIP.	p. 6
XKTR_PARAM_MIP_INTGAPREL	The relative integrality gap stop tolerance for MIP.	p. 6
XKTR_PARAM_MIP_KNAPSACK	Specifies rules for adding MIP knapsack cuts.	p. 15
XKTR_PARAM_MIP_LPALG	Specifies which algorithm to use for any linear programming (LP) subproblem solves that may occur in the MIP branch and bound procedure.	p. 16

XKTR_PARAM_MIP_MAXNODES	Specifies the maximum number of nodes explored.	p. 16
XKTR_PARAM_MIP_MAXSOLVES	Specifies the maximum number of subproblem solves allowed (0 means no limit).	p. 16
XKTR_PARAM_MIP_METHOD	Specifies which MIP method to use.	p. 17
XKTR_PARAM_MIP_OUTINTERVAL	Specifies node printing interval for XKTR_PARAM_MIP_OUTLEVEL when XKTR_PARAM_MIP_OUTLEVEL > 0.	p. 17
XKTR_PARAM_MIP_OUTLEVEL	Specifies how much MIP information to print.	p. 17
XKTR_PARAM_MIP_PSEUDOINIT	Specifies the method used to initialize pseudo-costs corresponding to variables that have not yet been branched on in the MIP method.	p. 17
XKTR_PARAM_MIP_ROOTALG	Specifies which algorithm to use for the root node solve in MIP (same options as XKTR_PARAM_ALGORITHM user option).	p. 18
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XKTR_PARAM_MIP_STRONG_LEVEL	Specifies the maximum number of tree levels on which to perform MIP strong branching.	p. 19
XKTR_PARAM_MIP_STRONG_MAXIT	Specifies the maximum number of iterations to allow for MIP strong branching solves.	p. 19
XKTR_PARAM_MIP_TERMINATE	Specifies conditions for terminating the MIP algorithm.	p. 19
XKTR_PARAM_OBJRANGE	Specifies the extreme limits of the objective function for purposes of determining unboundedness.	p. 6
XKTR_PARAM_OPTTOL	Specifies the final relative stopping tolerance for the KKT (optimality) error.	p. 6
XKTR_PARAM_OPTTOLABS	Specifies the final absolute stopping tolerance for the KKT (optimality) error.	p. 6
XKTR_PARAM_OUTLEV	Controls the level of output produced by KNITRO.	p. 19
XKTR_PARAM_PRESOLVE	Determine whether or not to use the KNITRO presolver to try to simplify the model by removing variables or constraints. Specifies conditions for terminating the MIP algorithm.	p. 20
XKTR_PARAM_PRESOLVE_TOL	Determines the tolerance used by the KNITRO presolver to remove variables and constraints from the model.	p. 7
XKTR_PARAM_SCALE	Performs a scaling of the objective and constraint functions based on their values at the initial point.	p. 20
XKTR_PARAM_SOC	Specifies whether or not to try second order corrections (SOC).	p. 20
XKTR_PARAM_XTOL	The optimization process will terminate if the relative change in all components of the solution point estimate is less than xtol.	p. 7

## 1.1 Double control parameters

These double control parameters can be set using XSLPsetdblcontrol using the Xpress-SLP API, XNLPsetsolverdoublecontrol using the XNLP API and setparam in Mosel using module mmxnlp.

---

### XKTR\_PARAM\_BAR\_FEASMODETOL

---

<b>Description</b>	Specifies the tolerance in equation that determines whether KNITRO will force subsequent iterates to remain feasible.
<b>Type</b>	Double
<b>Note</b>	The tolerance applies to all inequality constraints in the problem. This option only has an effect if option <code>XKTR_PARAM_BAR_FEASIBLE = stay</code> or <code>XKTR_PARAM_BAR_FEASIBLE = get_stay</code> .
<b>Default value</b>	1.0e-4

---

### XKTR\_PARAM\_BAR\_INITMU

---

<b>Description</b>	Specifies the initial value for the barrier parameter : $\mu$ used with the barrier algorithms. This option has no effect on the Active Set algorithm.
<b>Type</b>	Double
<b>Default value</b>	1.0e-1

---

### XKTR\_PARAM\_DELTA

---

<b>Description</b>	Specifies the initial trust region radius scaling factor used to determine the initial trust region size.
<b>Type</b>	Double
<b>Default value</b>	1.0e0

---

### XKTR\_PARAM\_DELTA

---

<b>Description</b>	Specifies the initial trust region radius scaling factor used to determine the initial trust region size.
<b>Type</b>	Double
<b>Default value</b>	1.0e0

---

## XKTR\_PARAM\_FEASTOL

---

<b>Description</b>	Specifies the final relative stopping tolerance for the feasibility error.
<b>Type</b>	Double
<b>Note</b>	Smaller values of feastol result in a higher degree of accuracy in the solution with respect to feasibility.
<b>Default value</b>	1.0e-6

---

## XKTR\_PARAM\_FEASTOLABS

---

<b>Description</b>	Specifies the final absolute stopping tolerance for the feasibility error.
<b>Type</b>	Double
<b>Note</b>	Smaller values of feastol_abs result in a higher degree of accuracy in the solution with respect to feasibility.
<b>Default value</b>	0.0e0

---

## XKTR\_PARAM\_INFEASTOL

---

<b>Description</b>	Specifies the (relative) tolerance used for declaring infeasibility of a model.
<b>Type</b>	Double
<b>Note</b>	Smaller values of infeastol make it more difficult to satisfy the conditions KNITRO uses for detecting infeasible models. If you believe KNITRO incorrectly declares a model to be infeasible, then you should try a smaller value for infeastol.
<b>Default value</b>	1.0e-8

---

## XKTR\_PARAM\_MIP\_INTEGERTOL

---

<b>Description</b>	This value specifies the threshold for deciding whether or not a variable is determined to be an integer.
<b>Type</b>	Double
<b>Default value</b>	1.0e-8

## XKTR\_PARAM\_MIP\_INTGAPABS

---

**Description** The absolute integrality gap stop tolerance for MIP.

**Type** Double

**Default value** 1.0e-6

---

## XKTR\_PARAM\_MIP\_INTGAPREL

---

**Description** The relative integrality gap stop tolerance for MIP.

**Type** Double

**Default value** 1.0e-6

---

## XKTR\_PARAM\_OBJRANGE

---

**Description** Specifies the extreme limits of the objective function for purposes of determining unboundedness.

**Type** Double

**Note** If the magnitude of the objective function becomes greater than objrange for a feasible iterate, then the problem is determined to be unbounded and KNITRO proceeds no further.

**Default value** 1.0e20

---

## XKTR\_PARAM\_OPTTOL

---

**Description** Specifies the final relative stopping tolerance for the KKT (optimality) error.

**Type** Double

**Note** Smaller values of opttol result in a higher degree of accuracy in the solution with respect to optimality.

**Default value** 1.0e-6

---

## XKTR\_PARAM\_OPTTOLABS

---

**Description** Specifies the final absolute stopping tolerance for the KKT (optimality) error.

<b>Type</b>	Double
<b>Note</b>	Smaller values of opttol_abs result in a higher degree of accuracy in the solution with respect to optimality.
<b>Default value</b>	0.0e0

---

## XKTR\_PARAM\_PRESOLVE\_TOL

---

<b>Description</b>	Determines the tolerance used by the KNITRO presolver to remove variables and constraints from the model.
<b>Type</b>	Double
<b>Note</b>	If you believe the KNITRO presolver is incorrectly modifying the model, use a smaller value for this tolerance (or turn the presolver off).
<b>Default value</b>	1.0e-6

---

## XKTR\_PARAM\_XTOL

---

<b>Description</b>	The optimization process will terminate if the relative change in all components of the solution point estimate is less than xtol.
<b>Type</b>	Double
<b>Note</b>	If using the Interior/Direct or Interior/CG algorithm and the barrier parameter is still large, KNITRO will first try decreasing the barrier parameter before terminating.
<b>Default value</b>	1.0e-15

## 1.2 Integer control parameters

These integer control parameters can be set using XSLPsetintcontrol using the Xpress-SLP API, XNLPsetsolverintcontrol using the XNLP API and setparam in Mosel using module mmxnlp.

---

### XKTR\_PARAM\_ALGORITHM

---

<b>Description</b>	Indicates which algorithm to use to solve the problem												
<b>Type</b>	Integer												
<b>Values</b>	<table><tr><td>0</td><td>(auto) let KNITRO automatically choose an algorithm, based on the problem characteristics.</td></tr><tr><td>1</td><td>(direct) use the Interior/Direct algorithm.</td></tr><tr><td>2</td><td>(cg) use the Interior/CG algorithm.</td></tr><tr><td>3</td><td>(active) use the Active Set algorithm.</td></tr><tr><td>4</td><td>(sqp) use the SQP algorithm.</td></tr><tr><td>5</td><td>(multi) run all algorithms, perhaps in parallel.</td></tr></table>	0	(auto) let KNITRO automatically choose an algorithm, based on the problem characteristics.	1	(direct) use the Interior/Direct algorithm.	2	(cg) use the Interior/CG algorithm.	3	(active) use the Active Set algorithm.	4	(sqp) use the SQP algorithm.	5	(multi) run all algorithms, perhaps in parallel.
0	(auto) let KNITRO automatically choose an algorithm, based on the problem characteristics.												
1	(direct) use the Interior/Direct algorithm.												
2	(cg) use the Interior/CG algorithm.												
3	(active) use the Active Set algorithm.												
4	(sqp) use the SQP algorithm.												
5	(multi) run all algorithms, perhaps in parallel.												
<b>Default value</b>	0												

---

### XKTR\_PARAM\_BAR\_DIRECTINTERVAL

---

<b>Description</b>	Controls the maximum number of consecutive conjugate gradient (CG) steps before KNITRO will try to enforce that a step is taken using direct linear algebra.
<b>Type</b>	Integer
<b>Note</b>	This option is only valid for the Interior/Direct algorithm and may be useful on problems where KNITRO appears to be taking lots of conjugate gradient steps. Setting bar_directinterval to 0 will try to enforce that only direct steps are taken which may produce better results on some problems.
<b>Default value</b>	10

---

### XKTR\_PARAM\_BAR\_FEASIBLE

---

<b>Description</b>	Specifies whether special emphasis is placed on getting and staying feasible in the interior-point algorithms.								
<b>Type</b>	Integer								
<b>Values</b>	<table><tr><td>0</td><td>(no) No special emphasis on feasibility.</td></tr><tr><td>1</td><td>(stay) Iterates must satisfy inequality constraints once they become sufficiently feasible.</td></tr><tr><td>2</td><td>(get) Special emphasis is placed on getting feasible before trying to optimize.</td></tr><tr><td>3</td><td>(get_stay) Implement both options 1 and 2 above.</td></tr></table>	0	(no) No special emphasis on feasibility.	1	(stay) Iterates must satisfy inequality constraints once they become sufficiently feasible.	2	(get) Special emphasis is placed on getting feasible before trying to optimize.	3	(get_stay) Implement both options 1 and 2 above.
0	(no) No special emphasis on feasibility.								
1	(stay) Iterates must satisfy inequality constraints once they become sufficiently feasible.								
2	(get) Special emphasis is placed on getting feasible before trying to optimize.								
3	(get_stay) Implement both options 1 and 2 above.								

<b>Note</b>	This option can only be used with the Interior/Direct and Interior/CG algorithms. If bar_feasible = stay or bar_feasible = get_stay, this will activate the feasible version of KNITRO. The feasible version of KNITRO will force iterates to strictly satisfy inequalities, but does not require satisfaction of equality constraints at intermediate iterates. This option and the honorbnds option may be useful in applications where functions are undefined outside the region defined by inequalities. The initial point must satisfy inequalities to a sufficient degree; if not, KNITRO may generate infeasible iterates and does not switch to the feasible version until a sufficiently feasible point is found. Sufficient satisfaction occurs at a point $x$ if it is true for all inequalities that $c_l + tol \leq c(x) \leq c_u - tol$ . The constant tol is determined by the option bar_feasmodetol. If bar_feasible = get or bar_feasible = get_stay, KNITRO will place special emphasis on first trying to get feasible before trying to optimize.
<b>Default value</b>	0

---

## XKTR\_PARAM\_BAR\_INITPT

---

<b>Description</b>	Indicates whether an initial point strategy is used with barrier algorithms.
<b>Type</b>	Integer
<b>Values</b>	0 (auto) Let KNITRO automatically choose the strategy. 1 (yes) Shift the initial slacks and multipliers to improve barrier algorithm performance. 2 (no) Do no alter the initial slacks and multipliers.
<b>Note</b>	This option has no effect on the Active Set algorithm.
<b>Default value</b>	0

---

## XKTR\_PARAM\_BAR\_MAXBACKTRACK

---

<b>Description</b>	Indicates the maximum allowable number of backtracks during the linesearch of the Interior/Direct algorithm before reverting to a CG step.
<b>Type</b>	Integer
<b>Note</b>	Increasing this value will make the Interior/Direct algorithm less likely to take CG steps. If the Interior/Direct algorithm is taking a large number of CG steps (as indicated by a positive value for 'Gits' in the output), this may improve performance. This option has no effect on the Active Set algorithm.
<b>Default value</b>	3

---

## XKTR\_PARAM\_BAR\_MAXCROSSIT

---

<b>Description</b>	Specifies the maximum number of crossover iterations before termination.
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<b>Type</b>	Integer
<b>Note</b>	If the value is positive and the algorithm in operation is Interior/Direct or Interior/CG, then KNITRO will crossover to the Active Set algorithm near the solution. The Active Set algorithm will then perform at most bar_maxcrossover iterations to get a more exact solution. If the value is 0, no Active Set crossover occurs and the interior-point solution is the final result. If Active Set crossover is unable to improve the approximate interior-point solution, then KNITRO will restore the interior-point solution. In some cases (especially on large-scale problems or difficult degenerate problems) the cost of the crossover procedure may be significant - for this reason, crossover is disabled by default. Enabling crossover generally provides a more accurate solution than Interior/Direct or Interior/CG.
<b>Default value</b>	0

---

## XKTR\_PARAM\_BAR\_MAXREFACTOR

---

<b>Description</b>	Indicates the maximum number of refactorizations of the KKT system per iteration of the Interior/Direct algorithm before reverting to a CG step.
<b>Type</b>	Integer
<b>Note</b>	These refactorizations are performed if negative curvature is detected in the model. Rather than reverting to a CG step, the Hessian matrix is modified in an attempt to make the subproblem convex and then the KKT system is refactorized. Increasing this value will make the Interior/Direct algorithm less likely to take CG steps. If the Interior/Direct algorithm is taking a large number of CG steps (as indicated by a positive value for "CGits" in the output), this may improve performance. This option has no effect on the Active Set algorithm.
<b>Default value</b>	-1

---

## XKTR\_PARAM\_BAR\_MURULE

---

<b>Description</b>	Indicates which strategy to use for modifying the barrier parameter mu in the barrier algorithms.
<b>Type</b>	Integer

<b>Values</b>	0	(auto) Let KNITRO automatically choose the strategy.
	1	(monotone) Monotonically decrease the barrier parameter. Available for both barrier algorithms.
	2	(adaptive) Use an adaptive rule based on the complementarity gap to determine the value of the barrier parameter. Available for both barrier algorithms.
	3	(probing) Use a probing (affine-scaling) step to dynamically determine the barrier parameter. Available only for the Interior/Direct algorithm.
	4	(dampmpc) Use a Mehrotra predictor-corrector type rule to determine the barrier parameter, with safeguards on the corrector step. Available only for the Interior/Direct algorithm.
	5	(fullmpc) Use a Mehrotra predictor-corrector type rule to determine the barrier parameter, without safeguards on the corrector step. Available only for the Interior/Direct algorithm.
	6	(quality) Minimize a quality function at each iteration to determine the barrier parameter. Available only for the Interior/Direct algorithm.
<b>Note</b>		Not all strategies are available for both barrier algorithms. This option has no effect on the Active Set algorithm.
<b>Default value</b>	0	

---

## XKTR\_PARAM\_BAR\_PENCONS

---

<b>Description</b>	Indicates whether a penalty approach is applied to the constraints.	
<b>Type</b>	Integer	
<b>Values</b>	0	(auto) Let KNITRO automatically choose the strategy.
	1	(none) No constraints are penalized.
	2	(all) A penalty approach is applied to all general constraints.
<b>Note</b>		Using a penalty approach may be helpful when the problem has degenerate or difficult constraints. It may also help to more quickly identify infeasible problems, or achieve feasibility in problems with difficult constraints. This option has no effect on the Active Set algorithm.
<b>Default value</b>	0	

---

## XKTR\_PARAM\_BAR\_PENRULE

---

<b>Description</b>	Indicates which penalty parameter strategy to use for determining whether or not to accept a trial iterate.	
<b>Type</b>	Integer	
<b>Values</b>	0	(auto) Let KNITRO automatically choose the strategy.
	1	(single) Use a single penalty parameter in the merit function to weight feasibility versus optimality.
	2	(flex) Use a more tolerant and flexible step acceptance procedure based on a range of penalty parameter values.

**Note** This option has no effect on the Active Set algorithm.

**Default value** 0

---

## XKTR\_PARAM\_BAR\_SWITCHRULE

---

**Description** Indicates whether or not the barrier algorithms will allow switching from an optimality phase to a pure feasibility phase.

**Type** Integer

**Values**

0	(auto) Let KNITRO determine the switching procedure.
1	(never) Never switch to feasibility phase.
2	(level1) Allow switches to feasibility phase.
3	(level2) Use a more aggressive switching rule.

**Note** This option has no effect on the Active Set algorithm.

**Default value** 0

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## XKTR\_PARAM\_GRADOPT

---

**Description** Specifies how to compute the gradients of the objective and constraint functions.

**Type** Integer

**Values**

1	(exact) User provides a routine for computing the exact gradients.
2	(forward) KNITRO computes gradients by forward finite-differences.
3	(central) KNITRO computes gradients by central finite differences.

**Note** It is highly recommended to provide exact gradients if at all possible as this greatly impacts the performance of the code.

**Default value** 1

---

## XKTR\_PARAM\_HESSOPT

---

**Description** Specifies how to compute the (approximate) Hessian of the Lagrangian.

**Type** Integer

**Values**

1	(exact) User provides a routine for computing the exact Hessian.
2	(bfgs) KNITRO computes a (dense) quasi-Newton BFGS Hessian.
3	(sr1) KNITRO computes a (dense) quasi-Newton SR1 Hessian.
4	(finite_diff) KNITRO computes Hessian-vector products using finite-differences.
5	(product) User provides a routine to compute the Hessian-vector products.
6	(lbfsgs) KNITRO computes a limited-memory quasi-Newton BFGS Hessian (its size is determined by the option lmsize).

<b>Note</b>	Options hessopt = 4 and hessopt = 5 are not available with the Interior/Direct algorithm. KNITRO usually performs best when the user provides exact Hessians (hessopt = 1) or exact Hessian-vector products (hessopt = 5). If neither can be provided but exact gradients are available (i.e., gradopt = 1), then hessopt = 4 is recommended. This option is comparable in terms of robustness to the exact Hessian option and typically not much slower in terms of time, provided that gradient evaluations are not a dominant cost. If exact gradients cannot be provided, then one of the quasi-Newton options is preferred. Options hessopt = 2 and hessopt = 3 are only recommended for small problems ( $n \leq 1000$ ) since they require working with a dense Hessian approximation. Option hessopt = 6 should be used for large problems.
<b>Default value</b>	1

---

## XKTR\_PARAM\_HONORBNDs

---

<b>Description</b>	Indicates whether or not to enforce satisfaction of simple variable bounds throughout the optimization.
<b>Type</b>	Integer
<b>Values</b>	0 (no) KNITRO does not require that the bounds on the variables be satisfied at intermediate iterates. 1 (always) KNITRO enforces that the initial point and all subsequent solution estimates satisfy the bounds on the variables. 2 (initpt) KNITRO enforces that the initial point satisfies the bounds on the variables.
<b>Note</b>	This option and the bar_feasible option may be useful in applications where functions are undefined outside the region defined by inequalities.
<b>Default value</b>	2

---

## XKTR\_PARAM\_LMSIZE

---

<b>Description</b>	Specifies the number of limited memory pairs stored when approximating the Hessian using the limited-memory quasi-Newton BFGS option.
<b>Type</b>	Integer
<b>Note</b>	The value must be between 1 and 100 and is only used with <a href="#">XKTR_PARAM_HESSOPT</a> = 6. Larger values may give a more accurate, but more expensive, Hessian approximation. Smaller values may give a less accurate, but faster, Hessian approximation. When using the limited memory BFGS approach it is recommended to experiment with different values of this parameter.
<b>Default value</b>	10

## XKTR\_PARAM\_MAXCGIT

---

<b>Description</b>	Specifies the number of limited memory pairs stored when approximating the Hessian using the limited-memory quasi-Newton BFGS option.
<b>Type</b>	Integer
<b>Values</b>	0      Let KNITRO automatically choose a value based on the problem size. n      At most n>0 CG iterations may be performed during one minor iteration of KNITRO.
<b>Default value</b>	0

---

## XKTR\_PARAM\_MAXIT

---

<b>Description</b>	Specifies the maximum number of iterations before termination.
<b>Type</b>	Integer
<b>Values</b>	0      Let KNITRO automatically choose a value based on the problem type. Currently KNITRO sets this value to 10000 for LPs/NLPs and 3000 for MIP problems. n      At most n>0 iterations may be performed before terminating.
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_BRANCHRULE

---

<b>Description</b>	Specifies which branching rule to use for MIP branch and bound procedure.
<b>Type</b>	Integer
<b>Values</b>	0      (auto) Let KNITRO automatically choose the branching rule. 1      (most_frac) Use most fractional (most infeasible) branching. 2      (pseudocost) Use pseudo-cost branching. 3      (strong) Use strong branching (see options <a href="#">XKTR_PARAM_MIP_STRONG_CANDLIM</a> , <a href="#">XKTR_PARAM_MIP_STRONG_LEVEL</a> , <a href="#">XKTR_PARAM_MIP_STRONG_MAXIT</a> for further control of strong branching procedure).
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_GUB\_BRANCH

---

<b>Description</b>	Specifies whether or not to branch on generalized upper bounds (GUBs).
<b>Type</b>	Integer

<b>Values</b>	0	(no) Do not branch on GUBs.
	1	(yes) Allow branching on GUBs.
<b>Default value</b>	0	

---

## XKTR\_PARAM\_MIP\_HEURISTIC

---

<b>Description</b>	Specifies which MIP heuristic search approach to apply to try to find an initial integer feasible point.	
<b>Type</b>	Integer	
<b>Values</b>	0	(auto) Let KNITRO choose the heuristic to apply (if any).
	1	(none) No heuristic search applied.
	2	(feaspump) Apply feasibility pump heuristic.
	3	(mpec) Apply heuristic based on MPEC formulation.
<b>Note</b>	If a heuristic search procedure is enabled, it will run for at most mip_heuristic_maxit iterations, before starting the branch and bound procedure.	
<b>Default value</b>	0	

---

## XKTR\_PARAM\_MIP\_HEURISTIC\_MAXIT

---

<b>Description</b>	Specifies the maximum number of iterations to allow for MIP heuristic, if one is enabled.	
<b>Type</b>	Integer	
<b>Default value</b>	100	

---

## XKTR\_PARAM\_MIP\_IMPLICATNS

---

<b>Description</b>	Specifies whether or not to add constraints to the MIP derived from logical implications.	
<b>Type</b>	Integer	
<b>Values</b>	0	(no) Do not add constraints from logical implications.
	1	(yes) KNITRO adds constraints from logical implications.
<b>Default value</b>	1	

---

## XKTR\_PARAM\_MIP\_KNAPSACK

---

<b>Description</b>	Specifies rules for adding MIP knapsack cuts.
--------------------	---

<b>Type</b>	Integer
<b>Values</b>	0      (none) Do not add knapsack cuts. 1      (ineqs) Add cuts derived from inequalities only. 2      (ineqs_eqs) Add cuts derived from both inequalities and equalities.
<b>Default value</b>	1

---

## XKTR\_PARAM\_MIP\_LPALG

---

<b>Description</b>	Specifies which algorithm to use for any linear programming (LP) subproblem solves that may occur in the MIP branch and bound procedure.
<b>Type</b>	Integer
<b>Values</b>	0      (auto) Let KNITRO automatically choose an algorithm, based on the problem characteristics. 1      (direct) Use the Interior/Direct (barrier) algorithm. 2      (cg) Use the Interior/CG (barrier) algorithm. 3      (active) Use the Active Set (simplex) algorithm.
<b>Note</b>	LP subproblems may arise if the problem is a mixed integer linear program (MILP), or if using <code>XKTR_PARAM_MIP_METHOD</code> = HQG. (Nonlinear programming subproblems use the algorithm specified by the algorithm option.)
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_MAXNODES

---

<b>Description</b>	Specifies the maximum number of nodes explored.
<b>Type</b>	Integer
<b>Note</b>	Zero value means no limit.
<b>Default value</b>	100000

---

## XKTR\_PARAM\_MIP\_MAXSOLVES

---

<b>Description</b>	Specifies the maximum number of subproblem solves allowed (0 means no limit).
<b>Type</b>	Integer
<b>Default value</b>	200000

---

## XKTR\_PARAM\_MIP\_METHOD

---

<b>Description</b>	Specifies which MIP method to use.						
<b>Type</b>	Integer						
<b>Values</b>	<table><tr><td>0</td><td>(auto) Let KNITRO automatically choose the method.</td></tr><tr><td>1</td><td>(BB) Use the standard branch and bound method.</td></tr><tr><td>2</td><td>(HQA) Use the hybrid Quesada-Grossman method (for convex, nonlinear problems only).</td></tr></table>	0	(auto) Let KNITRO automatically choose the method.	1	(BB) Use the standard branch and bound method.	2	(HQA) Use the hybrid Quesada-Grossman method (for convex, nonlinear problems only).
0	(auto) Let KNITRO automatically choose the method.						
1	(BB) Use the standard branch and bound method.						
2	(HQA) Use the hybrid Quesada-Grossman method (for convex, nonlinear problems only).						
<b>Default value</b>	0						

---

## XKTR\_PARAM\_MIP\_OUTINTERVAL

---

<b>Description</b>	Specifies node printing interval for <a href="#">XKTR_PARAM_MIP_OUTLEVEL</a> when <a href="#">XKTR_PARAM_MIP_OUTLEVEL &gt; 0</a> .						
<b>Type</b>	Integer						
<b>Values</b>	<table><tr><td>0</td><td>Print output every node.</td></tr><tr><td>2</td><td>Print output every 2nd node.</td></tr><tr><td>N</td><td>Print output every Nth node.</td></tr></table>	0	Print output every node.	2	Print output every 2nd node.	N	Print output every Nth node.
0	Print output every node.						
2	Print output every 2nd node.						
N	Print output every Nth node.						
<b>Default value</b>	10						

---

## XKTR\_PARAM\_MIP\_OUTLEVEL

---

<b>Description</b>	Specifies how much MIP information to print.				
<b>Type</b>	Integer				
<b>Values</b>	<table><tr><td>0</td><td>(none) Do not print any MIP node information.</td></tr><tr><td>1</td><td>(iters) Print one line of output for every node.</td></tr></table>	0	(none) Do not print any MIP node information.	1	(iters) Print one line of output for every node.
0	(none) Do not print any MIP node information.				
1	(iters) Print one line of output for every node.				
<b>Default value</b>	1				

---

## XKTR\_PARAM\_MIP\_PSEUDOUNIT

---

<b>Description</b>	Specifies the method used to initialize pseudo-costs corresponding to variables that have not yet been branched on in the MIP method.
<b>Type</b>	Integer

<b>Values</b>	0	Let KNITRO automatically choose the method.
	1	Initialize using the average value of computed pseudo-costs.
	2	Initialize using strong branching.
<b>Default value</b>	0	

---

## XKTR\_PARAM\_MIP\_ROOTALG

---

<b>Description</b>	Specifies which algorithm to use for the root node solve in MIP (same options as <a href="#">XKTR_PARAM_ALGORITHM</a> user option).
<b>Type</b>	Integer
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_ROUNDING

---

<b>Description</b>	Specifies the MIP rounding rule to apply.
<b>Type</b>	Integer
<b>Values</b>	0      (auto) Let KNITRO choose the rounding rule. 1      (none) Do not round if a node is infeasible. 2      (heur_only) Round using a fast heuristic only. 3      (nlp_sometimes) Round and solve a subproblem if likely to succeed. 4      (nlp_always) Always round and solve a subproblem.
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_SELECTRULE

---

<b>Description</b>	Specifies the MIP select rule for choosing the next node in the branch and bound tree.
<b>Type</b>	Integer
<b>Values</b>	0      (auto) Let KNITRO choose the node selection rule. 1      (depth_first) Search the tree using a depth first procedure. 2      (best_bound) Select the node with the best relaxation bound. 3      (combo_1) Use depth first unless pruned, then best bound.
<b>Default value</b>	0

---

## XKTR\_PARAM\_MIP\_STRONG\_CANDLIM

---

<b>Description</b>	Specifies the maximum number of candidates to explore for MIP strong branching.
<b>Type</b>	Integer
<b>Default value</b>	10

---

---

## XKTR\_PARAM\_MIP\_STRONG\_LEVEL

---

<b>Description</b>	Specifies the maximum number of tree levels on which to perform MIP strong branching.
<b>Type</b>	Integer
<b>Default value</b>	10

---

---

## XKTR\_PARAM\_MIP\_STRONG\_MAXIT

---

<b>Description</b>	Specifies the maximum number of iterations to allow for MIP strong branching solves.
<b>Type</b>	Integer
<b>Default value</b>	1000

---

---

## XKTR\_PARAM\_MIP\_TERMINATE

---

<b>Description</b>	Specifies conditions for terminating the MIP algorithm.
<b>Type</b>	Integer
<b>Values</b>	0      (optimal) Terminate at optimum. 1      (feasible) Terminate at first integer feasible point.
<b>Default value</b>	0

---

---

## XKTR\_PARAM\_OUTLEV

---

<b>Description</b>	Controls the level of output produced by KNITRO.
<b>Type</b>	Integer

<b>Values</b>	0	(none) Printing of all output is suppressed.
	1	(summary) Print only summary information.
	2	(iter_10) Print basic information every 10 iterations.
	3	(iter) Print basic information at each iteration.
	4	(iter_verbose) Print basic information and the function count at each iteration.
	5	(iter_x) Print all the above, and the values of the solution vector x.
	6	(all) Print all the above, and the values of the constraints c at x and the Lagrange multipliers lambda.
<b>Default value</b>	2	

---

## XKTR\_PARAM\_PRESOLVE

---

<b>Description</b>	Determine whether or not to use the KNITRO presolver to try to simplify the model by removing variables or constraints. Specifies conditions for terminating the MIP algorithm.
<b>Type</b>	Integer
<b>Values</b>	0 (none) Do not use KNITRO presolver. 1 (basic) Use the KNITRO basic presolver.
<b>Default value</b>	1

---

## XKTR\_PARAM\_SCALE

---

<b>Description</b>	Performs a scaling of the objective and constraint functions based on their values at the initial point.
<b>Type</b>	Integer
<b>Values</b>	0 (no) No scaling is performed. 1 (yes) KNITRO is allowed to scale the objective function and constraints.
<b>Note</b>	If scaling is performed, all internal computations, including the stopping tests, are based on the scaled values.
<b>Default value</b>	1

---

## XKTR\_PARAM\_SOC

---

<b>Description</b>	Specifies whether or not to try second order corrections (SOC).
<b>Type</b>	Integer

<b>Values</b>	0	(no) No second order correction steps are attempted.
	1	(maybe) Second order correction steps may be attempted on some iterations.
	2	(yes) Second order correction steps are always attempted if the original step is rejected and there are nonlinear constraints.
<b>Note</b>	A second order correction may be beneficial for problems with highly nonlinear constraints.	
<b>Default value</b>	1	

# APPENDIX A

## Contacting FICO

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FICO provides clients with support and services for all our products. Refer to the following sections for more information.

### Product support

FICO offers technical support and services ranging from self-help tools to direct assistance with a FICO technical support engineer. Support is available to all clients who have purchased a FICO product and have an active support or maintenance contract. You can find support contact information on the Product Support home page ([www.fico.com/support](http://www.fico.com/support)).

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### Product documentation

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## Sales and maintenance

*USA, CANADA AND ALL AMERICAS*

*Email:* XpressSalesUS@fico.com

*WORLDWIDE*

*Email:* XpressSalesUK@fico.com

*Tel:* +44 207 940 8718

*Fax:* +44 870 420 3601

Xpress Optimization, FICO

FICO House

International Square

Starley Way

Birmingham B37 7GN

UK

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