



Improving Skipping Fault Correction Attacks on Randomized Dilithium via MILP

Haobo Ouyang Chaoran Wang Guowei Liu
Lixuan Wu Meiqin Wang Yanhong Fan

School of Cyber Science and Technology, Shandong University

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Core signing computation

$$\mathbf{z} = \mathbf{y} + c\mathbf{s}_1$$

- $\mathbf{y}$ is freshly randomized for every signature.
- c is a sparse challenge with exactly τ nonzero coefficients.
- $\mathbf{s}_1$ is secret, but each coefficient is small: $s[j] \in [-\eta, \eta]$.

Randomized signing does not fully preclude fault-based key recovery.

Relevant parameters ($N = 256$)

Level	ℓ	η	τ	ℓN
L2	4	2	39	1024
L3	5	4	49	1280
L5	7	2	60	1792

Key property: the coefficients of $\mathbf{s}_1$ lie in the bounded small-integer domain $[-\eta, \eta]$.

One Fault, One Equation

- 1 **Inject an instruction-skipping fault** while computing $\mathbf{z} = \mathbf{y} + c\mathbf{s}_1$.
- 2 **Use public verification as a correction oracle.** At the faulted coefficient, try $\alpha \in [-\beta, \beta]$ until the modified signature verifies.
- 3 **Record the recovered relation**

$$\langle c_i, \mathbf{s} \rangle = \alpha,$$

where c_i is a rotation of the public challenge and $\mathbf{s}$ is one N -coefficient component of $\mathbf{s}_1$.

Krahmer et al.: collect $N = 256$ equations for every component to obtain a full-rank system.

Why Full Rank Is Unnecessary

Baseline: full-rank equations

- $M = N = 256$ equations per component
- Collect a full-rank system for each component
- Total number of injected faults: ℓN
- Does not use $\mathbf{s} \in [-\eta, \eta]^N$ in recovery

Our approach: underdetermined + bounded

- Collect only $M < N$ equations
- Retain the coefficient bound explicitly
- Restrict recovery to the bounded small-integer domain
- Recover with an MILP solver

Question. How small can M be while

$$\mathcal{S} = \{\mathbf{s} \in \mathbb{Z}^N : C\mathbf{s} = \mathbf{x}, \|\mathbf{s}\|_{\infty} \leq \eta\}$$

still contains exactly one element with high probability?

Online: collect corrected faults

- Each usable fault returns (c_i, x_i) .
- Every row has τ entries in $\{\pm 1\}$.
- Stack M samples into $C \in \{-1, 0, 1\}^{M \times N}$ and $\mathbf{x} \in \mathbb{Z}^M$.

Offline: solve a bounded system

$$C\mathbf{s} = \mathbf{x}, \quad \mathbf{s} \in [-\eta, \eta]^N.$$

- The equation system is underdetermined when $M < N$.
- The coefficient-wise bound limits the search domain.
- The bounded feasible set can still contain exactly one element.

The online phase collects equations; the offline phase invokes the MILP solver.

Step 1: coverage

- Let H_j count rows involving coefficient j .
- If $H_j = 0$, coefficient j is unconstrained; uniqueness is impossible.
- Under the random-row model,

$$\Pr[\forall j: H_j \geq 1] \geq 1 - N \left(1 - \frac{\tau}{N}\right)^M.$$

Step 2: uniqueness

- Two bounded solutions differ by a nonzero $\mathbf{d} \in [-2\eta, 2\eta]^N$.
- Non-uniqueness is equivalent to $\mathbf{C}\mathbf{d} = 0$ for an admissible $\mathbf{d}$.
- A first-moment bound gives

$$\Pr[\text{non-unique}] \leq ((2\eta + 1)^N - 1) p_{\text{gauss}}^M.$$

Full coverage is necessary; Theorem 1 bounds the probability of a second bounded solution.

Sufficient Fault Thresholds

For target failure probability $2^{-\lambda}$, it suffices to choose

$$M \geq \frac{\log((2\eta + 1)^N - 1) + \lambda}{-\log p_{\text{gauss}}}.$$

Here

$$p_{\text{gauss}} = \frac{1}{\sqrt{2\pi} \sigma_d}, \quad \sigma_d^2 = \frac{2\tau\eta(\eta + 1)}{3}.$$

This sufficient condition is derived under the random-row model and is not intended to be tight.

Level	p_{gauss}	$M_{\min}$	$M_{\min}/N$
L2	0.03194	122	47.7%
L3	0.01561	137	53.5%
L5	0.02575	114	44.5%

Values are per polynomial component and target at least 99% uniqueness probability.

Lemma 1

- Each row contains τ nonzero positions.
- Position j is absent from all M rows with probability $(1 - \frac{\tau}{N})^M$.
- A union bound over all N positions gives

$$\Pr[\exists j : H_j = 0] \leq N \left(1 - \frac{\tau}{N}\right)^M.$$

Dilithium-L2 example

$$N = 256, \quad \tau = 39, \quad M = 128.$$

$$256 \left(1 - \frac{39}{256}\right)^{128} \approx 1.7 \times 10^{-7}.$$

Therefore, at $M = N/2$, all secret coefficients are covered with high probability.

Coverage is necessary but does not by itself guarantee a unique bounded solution.

MILP Recovery Model

Variables

$$s[j] \in \mathbb{Z}, \quad -\eta \leq s[j] \leq \eta$$

$N = 256$ bounded integers

Constraints

$$C_i \mathbf{s} = x_i, \quad i = 0, \dots, M - 1$$

$M < N$ sparse equalities

Solver setup

$$\min f(\mathbf{s})$$

MILP solver (Gurobi)
objective function $f(\mathbf{s})$

Adaptive retry strategy

- 1 Start below N with M_{try} collected equations.
- 2 Solve one component and check the candidate by reconstructing signatures and running public verification.
- 3 If the check fails or solving times out, collect four more equations and retry.

Plain and Shuffling Settings

Plain setting

- Fault location is controlled and known.
- Repeatedly target a fixed coefficient in each component.
- Append each corrected equation to that component's set.
- Invoke MILP after M_{try} samples; retry adaptively.

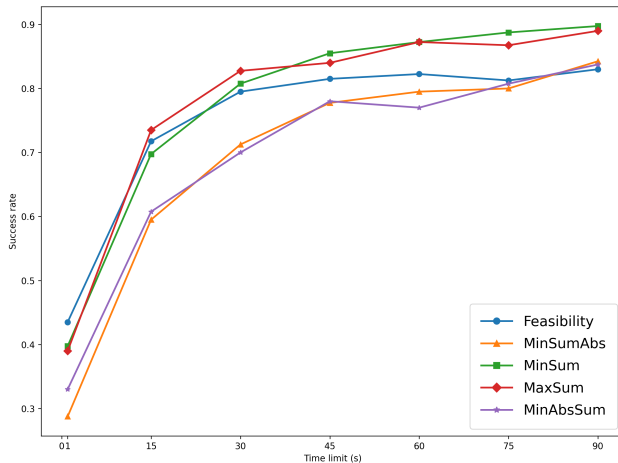
The target position is known

Shuffling setting

- Addition order is randomized; the fault location is unknown.
- Search all ℓN candidate locations during correction.
- Route a recovered equation to the corresponding component.
- Stop solving components once their candidates verify.

The fault location must be resolved by correction search

Objective and Time Limit



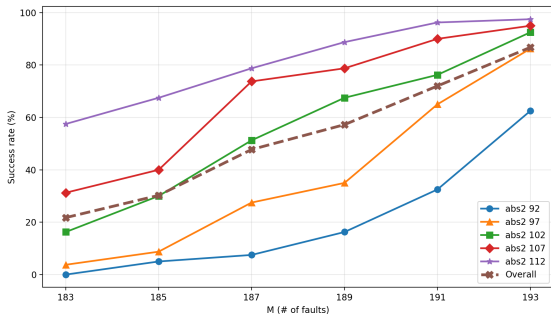
Setup

- 11th Gen Intel i7 at 2.5 GHz
- Gurobi MILP solver
- Dilithium implementation
- L2; $M = 193$; 400 components

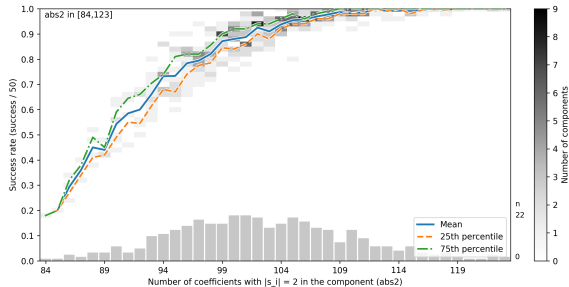
Observation

- Feasibility leads for short budgets.
- MinSum/MaxSum lead after 30 seconds.
- Curves largely plateau by $T = 60$ s.

Recovery Success Factors



At fixed abs2, success rises with M .



At fixed $M = 193$, components with larger abs2 are easier.

Explanation: larger $\|s\|_2^2$ lowers the single-row collision probability, making the system more constrained.

Fault Reduction: 13–26%

Plain setting

Level	Baseline	Ours	Reduction
L2	1024	759	25.9%
L3	1280	1073	16.2%
L5	1792	1334	25.6%

200 independent trials; 60-second MILP budget.

Shuffling setting

Level	Baseline	Ours	Reduction
L2	1157	861	25.6%
L3	1374	1189	13.5%
L5	2075	1529	26.3%

10 independent trials; exhaustive location search.

Under shuffling, each usable equation requires correction search over all ℓN possible coefficient positions.

$M_{\min} = 122$ (L2 sufficient uniqueness bound) vs. $M = 193$
(97.5% at $\text{abs2} = 112$)

What the theorem controls

- An idealized independent random-row model
- Existence of a unique bounded solution
- A conservative first-moment/union bound

What the experiment adds

- Generic branch-and-bound search
- A finite 60-second time budget
- Private-key feature `abs2`
- Matrix rows obtained from rotated challenge vectors

Interpretation: uniqueness is necessary for recovery, but not sufficient for fast recovery by a generic solver.

Sign-then-verify

- Verification runs inside the signing device.
- A signature that fails internal verification is not returned.
- The attacker cannot obtain the faulty signature required by EQFROMSIG.
- Public verification can no longer be used as the correction oracle.

Double computation

- The signing computation is performed twice.
- Results are compared before releasing the signature.
- A successful attack requires compatible faults in both executions.
- Such faults typically have to affect the same coefficient position.

The paper does not evaluate the attack under these two countermeasures; quantitative analysis is left for future work.

Key Takeaways

- 1 **Theory:** coverage plus bounded-solution uniqueness gives $(M_{\min})_{L2,L3,L5} = (122, 137, 114)$ per component for a 99% target.
- 2 **Method:** MILP makes the coefficient bound explicit and supports adaptive online collection with offline retry.
- 3 **Attack settings:** the adaptive MILP attack is described for both plain and shuffling settings.
- 4 **Experimental result:** the measured number of injected faults drops by 13–26% across all three Dilithium security levels.

Collecting a full-rank system of equations is not required for key recovery.

Questions & Discussion

Haobo Ouyang

hbouyang@mail.sdu.edu.cn

Code: <https://anonymous.4open.science/r/sfcafDvM-H73B>