# ddml: Double/debiased machine learning in Stata

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Package website: https://statalasso.github.io/

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# Introduction

- A rich and growing literature exploits machine learning to facilitate causal inference.
- A central focus: high-dimensional controls and/or instruments, which can arise if
  - ► we observe many controls/instruments
  - controls/instruments enter through an unknown function
- Belloni, Chernozhukov, and Hansen (2014) and Belloni et al. (2012) propose estimators *relying on the Lasso* that allow for high-dimensional controls/instruments.
  - $\Rightarrow$  Available via pdslasso in Stata (Ahrens, Hansen, and Schaffer, 2020)

# Introduction

#### What if we don't want to use the lasso?

- The Lasso might not be the best-performing machine learner for a particular problem.
- ► The Lasso relies on the *approximate sparsity assumption*, which might not be appropriate in some settings.

Chernozhukov et al. (2018) propose *Double/Debiased Machine Learning* (DDML) which allow to exploit machine learners other than the Lasso.

#### Our contribution:

- ▶ We introduce ddml, which implements DDML for Stata.
- We provide simulation evidence on the finite sample performance of DDML.
- Our recommendation is to use DDML in combination with stacking.

Motivating example. The partial linear model:



How do we account for confounding factors  $\mathbf{x}_i$ ? — The standard approach is to assume linearity  $g(\mathbf{x}_i) = \mathbf{x}'_i \beta$  and consider alternative combinations of controls.

#### Problems:

- Non-linearity & unknown interaction effects
- ► High-dimensionality: we might have "many" controls
- ► We don't know which controls to include

Motivating example. The partial linear model:



*Post-double selection* (Belloni, Chernozhukov, and Hansen, 2014) and *post-regularization* (Chernozhukov, Hansen, and Spindler, 2015) provide data-driven solutions for this setting.

Both "double" approaches rely on the *sparsity assumption* and use two auxiliary lasso regressions:  $y_i \rightsquigarrow \mathbf{x}_i$  and  $d_i \rightsquigarrow \mathbf{x}_i$ . Lasso PDS

Related approaches exist for *optimal IV* estimation (Belloni et al., 2012) and/or *IV with many controls* (Chernozhukov, Hansen, and Spindler, 2015).

These methods have been implemented for Stata in pdslasso (Ahrens, Hansen, and Schaffer, 2020), dsregress (StataCorp) and R (hdm; Chernozhukov, Hansen, and Spindler, 2016).

```
A quick example using AJR (2001):
```

```
. clear
```

```
. use https://statalasso.github.io/dta/AJR.dta
```

```
. pdslasso logpgp95 avexpr ///
(lat_abst edes1975 avelf temp* humid* steplow-oilres)
```

Example 1 (pdslasso) allows for high-dimensional controls.

*Example 2* (ivlasso) treats avexpr as endogenous and exploits logem4 as an instrument. (More details in the pds/ivlasso help file.)

There are **advantages** of relying on lasso:

- ▶ intuitive assumption of (approximate) sparsity
- computationally relatively cheap (due to plugin lasso penalty; no cross-validation needed)
- Linearity has its advantages (e.g. extension to fixed effects; Belloni et al., 2016)

But there are also drawbacks:

- What if the sparsity assumption is not plausible?
- There is a wide set of machine learners at disposable—Lasso might not be the best choice.
- Lasso requires careful feature engineering to deal with non-linearity & interaction effects.

## **Review of DDML**

The partial linear model:

$$y_i = \theta d_i + g(\mathbf{x}_i) + \varepsilon_i$$
$$d_i = m(\mathbf{x}_i) + v_i$$

*Naive idea:* We estimate conditional expectations  $\ell(\mathbf{x}_i) = E[y_i | \mathbf{x}_i]$  and  $m(\mathbf{x}_i) = E[d_i | \mathbf{x}_i]$  using ML and partial out the effect of  $\mathbf{x}_i$  (in the style of Frisch-Waugh-Lovell):

$$\hat{\theta}_{DDML} = \left(\frac{1}{n}\sum_{i}\hat{v}_{i}^{2}\right)^{-1}\frac{1}{n}\sum_{i}\hat{v}_{i}(y_{i}-\hat{\ell}),$$

where  $\hat{v}_i = d_i - \hat{m}_i$ .

# **Review of DDML**

Yet, this approach is flawed: The estimation error  $\ell(\mathbf{x}_i) - \hat{\ell}$  and  $v_i$  may be correlated due to **over-fitting**, leading to poor performance.

DDML, thus, relies on **cross-fitting**. Cross-fitting is sample splitting with swapped samples.

DDML for the partial linear model (DML 2)

We split the sample in K random folds of equal size denoted by  $I_k$ :

- For k = 1,..., K, estimate ℓ(x<sub>i</sub>) and m(x<sub>i</sub>) using sample I<sup>c</sup><sub>k</sub> and form out-of-sample predictions ℓ<sub>i</sub> and m̂<sub>i</sub> for all i in I<sub>k</sub>.
- $\blacktriangleright$  Construct estimator  $\hat{\theta}$  as

$$\left(\frac{1}{n}\sum_{i}\hat{v}_{i}^{2}\right)^{-1}\frac{1}{n}\sum_{i}\hat{v}_{i}(y_{i}-\hat{\ell}),$$

where  $\hat{v}_i = d_i - \hat{m}_i$ .  $\hat{m}_i$  and  $\hat{\ell}_i$  are the cross-fitted predicted values.

The DDML framework can be applied to other models (all implemented in ddml):

Interactive model

$$y_i = g(d_i, \mathbf{x}_i) + u_i \qquad E[u_i | \mathbf{x}_i, d_i] = 0$$
  
$$z_i = m(\mathbf{x}_i) + v_i \qquad E[u_i | \mathbf{x}_i] = 0$$

As in the Partial Linear Model, we are interested in the ATE, but do not assume that  $d_i$  (a binary treatment variable) and  $x_i$  are separable.

We estimate the conditional expectations  $E[y_i|\mathbf{x}_i, d_i = 0]$  and  $E[y_i|\mathbf{x}_i, d_i = 1]$  as well as  $E[d_i|\mathbf{x}_i]$  using a supervised machine learner.

The DDML framework can be applied to other models (all implemented in ddml):

Partial linear IV model

$$y_i = d_i\theta + g(\mathbf{x}_i) + u_i \qquad E[u_i|\mathbf{x}_i, z_i] = 0$$
  
$$z_i = m(\mathbf{x}_i) + v_i \qquad E[v_i|\mathbf{x}_i] = 0$$

where the aim is to estimate the average treatment effect  $\theta$  using observed instrument  $z_i$  in the presence of controls  $x_i$ . We estimate the conditional expectations  $E[y_i|x_i]$ ,  $E[d_i|x_i]$  and  $E[z_i|x_i]$  using a supervised machine learner.

The DDML framework can be applied to other models (all implemented in ddml):

High-dimensional IV model

$$y_i = d_i\theta + g(\mathbf{x}_i) + u_i$$
  
$$d_i = h(\mathbf{z}_i) + m(\mathbf{x}_i) + v_i$$

where the estimand of interest is  $\theta$ . The instruments and controls enter the model through unknown functions g(), h() and f().

We estimate the conditional expectations  $E[y_i|\mathbf{x}_i]$ ,  $E[\hat{d}_i|\mathbf{x}_i]$  and  $\hat{d}_i := E[d_i|\mathbf{z}_i, \mathbf{x}_i]$  using a supervised machine learner. The instrument is then formed as  $\hat{d}_i - \hat{E}[\hat{d}_i|\mathbf{x}_i]$  where  $\hat{E}[\hat{d}_i|\mathbf{x}_i]$  denotes the estimate of  $E[\hat{d}_i|\mathbf{x}_i]$ .

The DDML framework can be applied to other models (all implemented in ddml):

Interactive IV model

$$y_{i} = \mu(\mathbf{x}_{i}, \mathbf{z}_{i}) + u_{i} \qquad E[u_{i}|\mathbf{x}_{i}, z_{i}] = 0$$
  
$$d_{i} = m(z_{i}, \mathbf{x}_{i}) + v_{i} \qquad E[v_{i}|\mathbf{x}_{i}, z_{i}] = 0$$
  
$$z_{i} = p(\mathbf{x}_{i}) + \xi_{i} \qquad E[\xi_{i}|\mathbf{x}_{i}] = 0$$

where the aim is to estimate the local average treatment effect.

We estimate, using a supervised machine learner, the following conditional expectations:  $E[y_i | \mathbf{x}_i, z_i = 0]$  and  $E[y_i | \mathbf{x}_i, z_i = 1]$ ;  $E[D|\mathbf{x}_i, z_i = 0]$  and  $E[D|\mathbf{x}_i, z_i = 1]$ ;  $E[z_i | \mathbf{x}_i]$ .

## The ddml package

We introduce ddml for Stata:

- Compatible with various ML programs in Stata (e.g. lassopack, pylearn, randomforest).
  - $\rightarrow$  Any program with the classical "reg y x" syntax and post-estimation predict will work.
- ▶ Short (one-line) and flexible multi-line version
- ► 5 models supported: partial linear model, interactive model, interactive IV model, partial IV model, optimal IV.

# Stacking regression

#### Which machine learner should we use?

ddml supports a range of ML programs: pylearn, lassopack, randomforest. — Which one should we use?

We don't know whether we have a sparse or dense problem; linear or non-linear. We don't know whether, e.g., lasso or random forests will perform better.

Stacking, as implemented in pystacked, provides a solution: We use an 'optimal' combination of base learners.

#### qddml example: Partial linear model

qddml is the one-line ('quick') version of ddml and uses a syntax similar to pds/ivlasso.

. use https://statalasso.github.io/dta/AJR.dta, clear

- . global Y logpgp95
- . global X lat\_abst edes1975 avelf temp\* humid\* steplow-oilres
- . global D avexpr

. qddml \$Y \$D (\$X), model(partial) cmdopt(method(rf gradboost))
DML with Y=m0\_y and D=m0\_d1:

m0_y	Coef.	Std. Err.	z	P> z	[95% Conf.	Interval]
m0_d1	.3391897	.0621291	5.46	0.000	.217419	.4609604

#### Extended ddml syntax

**Step 1:** Initialise ddml and select model:

ddml init model [, kfolds(integer) reps(integer) ...] where model is either 'partial', 'iv', 'interactive', 'ivhd', 'late'.

**Step 2:** Add supervised ML programs for estimating conditional expectations:

```
ddml eq newvarname [, eqopt]: command depvar indepvars [,
cmdopt]
```

where *eq* selects the conditional expectations to be estimated. *command* is a ML program that supports the standard reg y x-type syntax. *cmdopt* are specific to that program.

Multiple estimation commands per equation are allowed.

#### Extended ddml syntax

Step 3: Cross-fitting

ddml crossfit

Step 4: Estimation of causal effects

ddml estimate [, robust ...]

#### Additional auxiliary commands:

ddml describe (describe current model set up),ddml save & ddml use (to import/save ddml objects), ddml extract (to retrieve objects), ddml export (export in csv format).

## Extended ddml syntax: Example

```
. global Y logpgp95
```

. global X lat\_abst edes1975 avelf temp\* humid\* steplow-oilres

```
. global D avexpr
```

```
. *** initialise ddml and select model;
```

```
. ddml init partial
```

```
. *** specify supervised machine learners for E[Y|X] ("yeq") and E[D|X] ("deq") . * y-equation:
```

```
. ddml yeq, gen(pyy): pystacked $Y $X, type(reg) method(rf gradboost) Equation successfully added.
```

```
. * d-equation:
```

. ddml deq, gen(pyd): pystacked \$D \$X, type(reg) method(rf gradboost) Equation successfully added.

# Extended ddml syntax: Example (cont'd.)

m0\_pyd

.3794184

. *** cross-fittin . ddml crossfit Model: partial Number of Y estima	ng and display mean ating equations: 1	-squared predic	ction e	)rror	
Cross-fitting equa	ation 1 2				
Mean-squared error Name	r for y X: Orthogonalized	Command	N	MSPE	
logpgp95	m0_pyy	pystacked	64	0.573751	
Mean-squared error Name	for D X: Orthogonalized	Command	N	MSPE	
avexpr	m0_pyd	pystacked	64	1.648804	
<ul> <li>. *** estimation of</li> <li>. ddml estimate</li> <li>DML with Y=m0 pvv</li> </ul>	of parameter of into and D=m0 pvd (N=):	erest			
m0_pyy	Coef. Std. Err	. z P>	z	[95% Conf. Interval]	_

6.67

0.000

.2678821

.0569073

.4909546

We demonstrate the use of ddml using the partially linear model by extending the analysis of 401(k) eligibility and total financial wealth of Poterba, Venti, and Wise (1995). The data consists of n = 9915 households from the 1991 SIPP.

**Simulation set-up:** We consider a *linear DGP* and a *non-linear DGP*, and compare performance of OLS, PDS-Lasso and various machine learners, including meta learners.

We would expect that stacking performs well under both settings, while linear approaches only perform well if the DGP is linear.

- 1. Use the full sample OLS estimate to obtain  $\hat{\tau}_{OLS}$ . Construct the partial residuals  $y_i^{(r)} = y_i \hat{\tau}_{OLS} d_i$ .
- 2. Fit a supervised learning estimator that aims to predict  $y_i^{(r)}$  with the controls  $x_i$  and  $d_i$  with  $x_i$ , respectively. Denote the fitted values as  $\tilde{g}$  and  $\tilde{h}$ . We either use
  - ► linear regression (*Linear DGP*)
  - gradient boosting (Non-linear DGP)
- 3. Sample from the empirical distribution of  $x_i$  by bootstrapping  $n_s$  observations from the original data. Denote the bootstrapped sample by  $\mathcal{D}_b$ .
- 4. Generate the treatment and outcome variable over the bootstrap sample:

$$\tilde{d}_{i}^{(b)} = \mathbb{1}\{\tilde{h}(x_{i}) + \nu_{i} \ge 0.5\}$$
 (1)

$$\tilde{y}_i^{(b)} = \tau_0 \tilde{d}_i^{(b)} + \tilde{g}(x_i) + \varepsilon_i, \forall i \in \mathcal{D}_b$$
(2)

where  $\nu_i \stackrel{iid}{\sim} \mathcal{N}(0, \kappa_1)$  and  $\varepsilon_i \stackrel{iid}{\sim} \mathcal{N}(0, \kappa_2)$ ,  $\tau_0 = 6,000$ .

We consider the following base learners:

- CV-Lasso with interactions and 2nd order polynomials
- CV-Ridge with interactions and 2nd order polynomials
- CV-Lasso with 10th order polynomials and no interactions
- CV-Ridge with 10th order polynomials and no interactions
- Random forest with low regularization: 8 predictors considered at each leaf split, no limit on the number of observation per node, bootstrap sample size of 70% (max\_features(8) min\_samples\_leaf(1) max\_samples(.7))
- Random forest with high regularization: 5 predictors considered at each leaf split, at least 10 observation per node, bootstrap sample size of 70% (max\_features(5) min\_samples\_leaf(10) max\_samples(.7))
- Gradient boosted trees with low regularization: 500 trees and a learnings rate of 0.01 (n\_estimators(500) learning\_rate(0.01))
- Gradient boosted trees with high regularization: 250 trees and a learnings rate of 0.01 (n\_estimators(250) learning\_rate(0.01))
- Feed-forward neural nets with 3 hidden layers of size 5 (hidden\_layer\_sizes(5 5 5))

Table: Average Stacking Weights

	Stacking		Single	-Best	
Panel (A): Linear DGP	Y X	D X	Y X	D X	
OLS	.646	.492	.813	.643	
Lasso with CV (2nd order poly)	.111	.158	.161	.267	
Ridge with CV (2nd order poly)	.063	.061	.018	.019	
Lasso with CV (10th order poly)	.032	.08	.003	.049	
Ridge with CV (10th order poly)	.03	.047	.005	.016	
Random forest (low regularization)	.013	.012	0	0	
Random forest (high regularization)	.017	.027	0	0	
Gradient boosting (low regularization)	.028	.043	0	.006	
Gradient boosting (high regularization)	.024	.074	0	.002	
Neural net	.036	.005	0	0	

Stacking assigns the highest weight to OLS if the DGP is linear...

Table: Average Stacking Weights

	Stacking		Single	e-Best	
	(1)	(2)	(3)	(4)	
Panel (B): Non-Linear DGP	Y X	D X	Y X	D X	
OLS	.037	.021	0	0	
Lasso with CV (2nd order poly)	.039	.067	.083	.149	
Ridge with CV (2nd order poly)	.177	.23	.12	.125	
Lasso with CV (10th order poly)	.052	.077	.088	.061	
Ridge with CV (10th order poly)	.078	.068	.019	.044	
Random forest (low regularization)	.041	.01	0	0	
Random forest (high regularization)	.028	.069	.001	0	
Gradient boosting (low regularization)	.517	.213	.678	.359	
Gradient boosting (high regularization)	.02	.239	.011	.262	
Neural net	.012	.005	0	0	

... and the highest weight to gradient boosting if *the DGP is non-linear*.

Table: Bias and Coverage Rates in the Linear and Non-Linear DGP

	$n_{s} = 9915$			$n_{s} = 99150$		
Panel (A): Linear DGP Full sample:	Bias	MAB	Rate	Bias	MAB	Rate
OLS	100.99	918.03	.95	-22.61	255.52	.94
PDS-Lasso	101.83	913.18	.95	-19.9	257.29	.94
DDML methods:						
Base learners						
OLS	105.07	906.96	.94	-23.05	256.51	.94
Lasso with CV (2nd order poly)	104.33	907.84	.94	-22.45	257.23	.94
Ridge with CV (2nd order poly)	103.22	898.56	.94	-23.27	255.54	.94
Lasso with CV (10th order poly)	49.56	1120.59	.93	37.98	260.53	.95
Ridge with CV (10th order poly)	1066	1342.38	.9	15.85	260.41	.95
Random forest (low regularization)	-59.63	1083.64	.91	-59.29	343.46	.86
Random forest (high regularization)	105.58	952.35	.94	-46.54	275.56	.91
Gradient boosting (low regularization)	53.97	930.93	.94	-41.84	252.14	.94
Gradient boosting (high regularization)	162.75	923.08	.95	48.31	259.12	.95
Neural net	-3594.99	5380.31	.17	-2165.41	3212.12	.16
Meta learners						
Stacking: NNLS	100.01	935.27	.94	-22.7	254.01	.94
Stacking: Single best	92.79	944.07	.95	-25.03	255.75	.94
Stacking: NNLS interactive	690.88	1245.02	.95	-51.29	359.4	.96
Short-stacking	100.51	912.88	.94	-24.04	252.17	.94
Single best	103.31	917.74	.94	-24.88	252.16	.94

Table: Bias and Coverage Rates in the Linear and Non-Linear DGP

	$n_{s} = 9915$			$n_{s} = 99150$			
Panel (B): Non-Linear DGP Full sample:	Bias	MAB	Rate	Bias	MAB	Rate	
OLS	-2496.16	2477.19	.63	-2658.04	2636.31	0	
PDS-Lasso	-2507.47	2489.77	.62	-2657.5	2635.94	0	
DDML methods:							
Base learners							
OLS	-2522.98	2540.36	.62	-2660.54	2640.98	0	
Lasso with CV (2nd order poly)	767.2	1078.29	.91	691.67	695.3	.64	
Ridge with CV (2nd order poly)	825.21	1091.19	.9	702.55	707.28	.64	
Lasso with CV (10th order poly)	-4214.09	1895.22	.92	-10.06	294.34	.94	
Ridge with CV (10th order poly)	-2123.59	2095.56	.91	4.42	288.37	.94	
Random forest (low regularization)	-104.54	1019.55	.92	-28.83	332.87	.87	
Random forest (high regularization)	-110.06	959.96	.95	-21.52	280.36	.94	
Gradient boosting (low regularization)	69.44	890.94	.95	7.28	263.62	.95	
Gradient boosting (high regularization)	213.04	895.47	.95	174.14	291.63	.93	
Neural net	-4706.76	5831.79	.17	-3216.85	3837.37	.15	
Meta learners							
Stacking: NNLS	-62.97	1068.87	.84	18.36	269.02	.95	
Stacking: Single best	-135.15	1035.41	.89	7.94	263.06	.95	
Stacking: NNLS interactive	118.88	1102.99	.95	10.27	280.24	.95	
Short-stacking	209.14	915.02	.95	13.67	261.73	.95	
Single best	125.64	914.56	.95	7.28	263.62	.95	

Wüthrich and Zhu (2021, henceforth WZ) consider two applications to demonstrate that PDS-Lasso suffers from a large finite sample bias and tends to underselect; again using the application of (Poterba, Venti, and Wise, 1995; Belloni et al., 2017).

They use two specifications:

- two-way interactions (TWI) (as in Chernozhukov and Hansen, 2004); p = 167
- quadratic splines & interactions (QSI) (as in Belloni et al., 2017); p = 272

WZ run their simulations on bootstrap samples of the data  $(n_b = \{200, 400, 800, 1600\})$  and calculate the bias as the mean difference to the full sample estimate (N = 9915).



(a) Bias (TWI specification)

(b) Bias (QSI specification)

*Notes:* The figures report the mean bias calculated as the mean difference to the full sample estimates. Following WZ, we draw 600 bootstrap samples of size  $n_b = \{200, 400, 600, 800, 1200, 1600\}$ . 'TWI' indicates that the predictors have been expanded by two-way interactions. 'QSI' refers to the quadratic spline & interactions specification of Belloni et al. (2017).

Figure: Replication of Figure 8 in WZ



(a) CV-Lasso

(b) CV-Ridge

Figure: Mean bias relative to full sample



(a) Boosted trees

(b) Stacking

Figure: Mean bias relative to full sample

**Main result:** The finite sample bias of stacking stabilizes for  $n_b > 600$ , but in contrast to OLS/PDS-Lasso, stacking doesn't assume linearity.

# Summary

- ddml implements Double/Debiased Machine Learning for Stata:
  - Compatible with various ML programs in Stata
  - Short (one-line) and flexible multi-line version
  - Uses Stacking Regression as the default machine learner; implemented via separate program pystacked
  - ► 5 models supported
- The advantage to pdslasso is that we can make use of almost any machine learner.
- But which machine learner should we use? We suggest Stacking as it combines multiple ML methods into one prediction.
- We are in the final phase of development; hopefully we can make ddml available soon (following your feedback)

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