

Figure S1

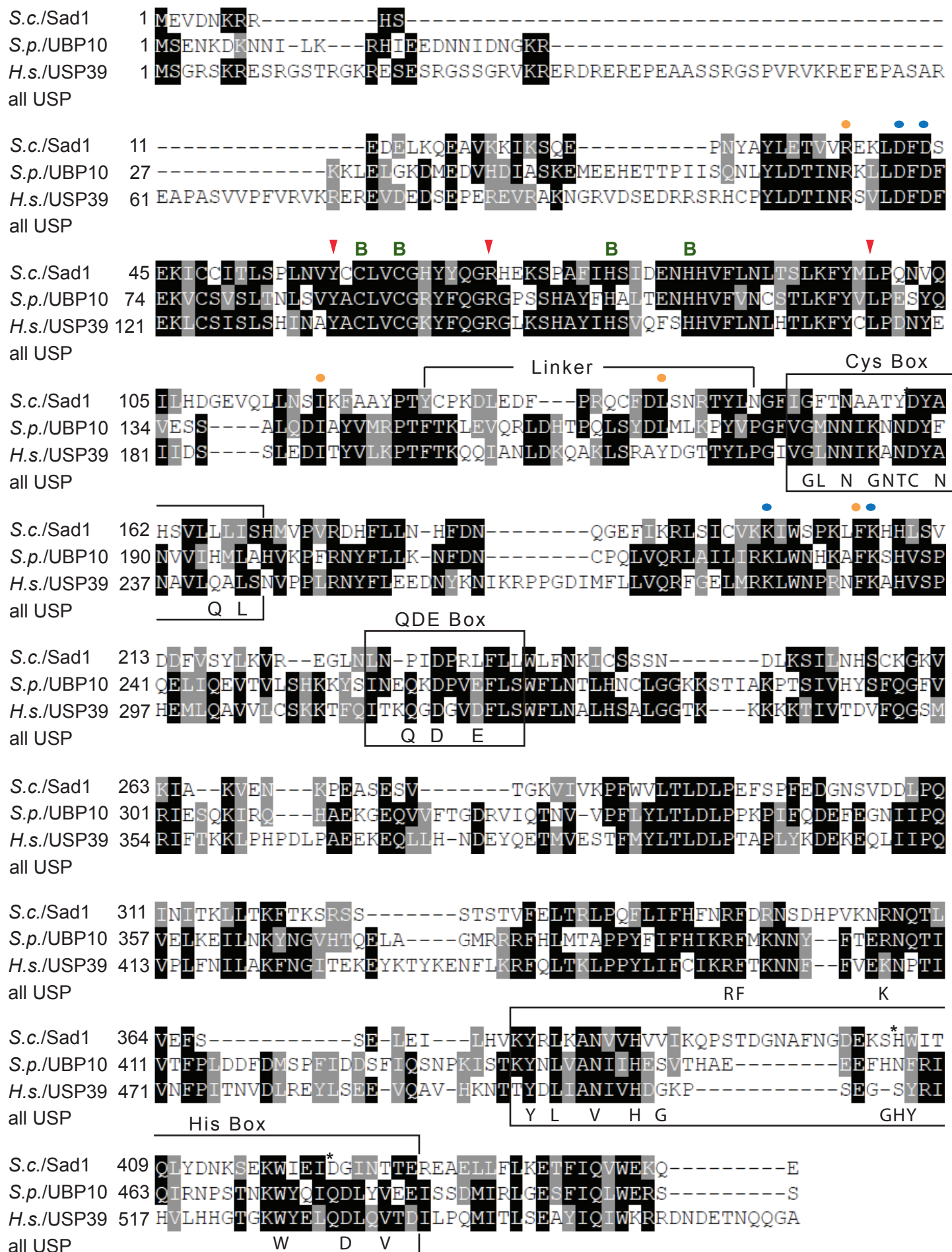
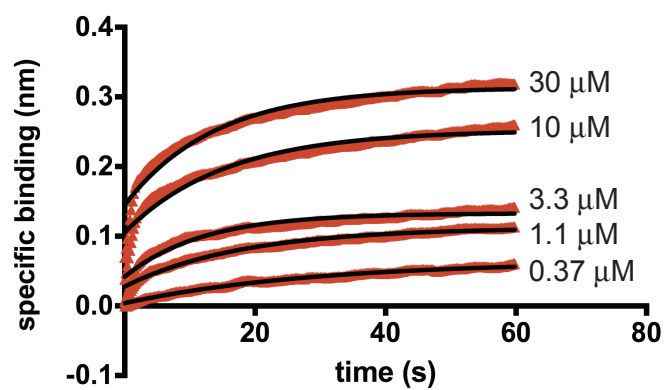


Figure S2

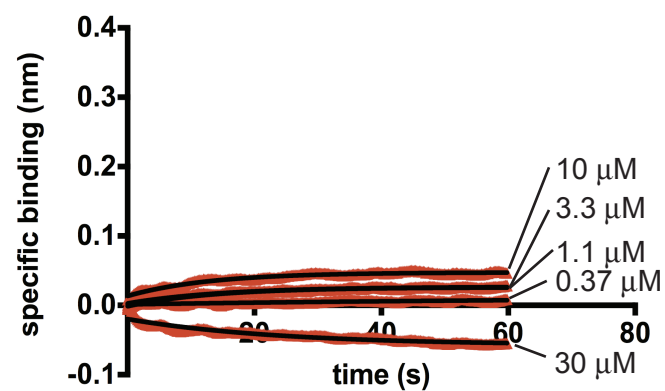
A

USP5 ZnF-UBP



B

Sad1



C

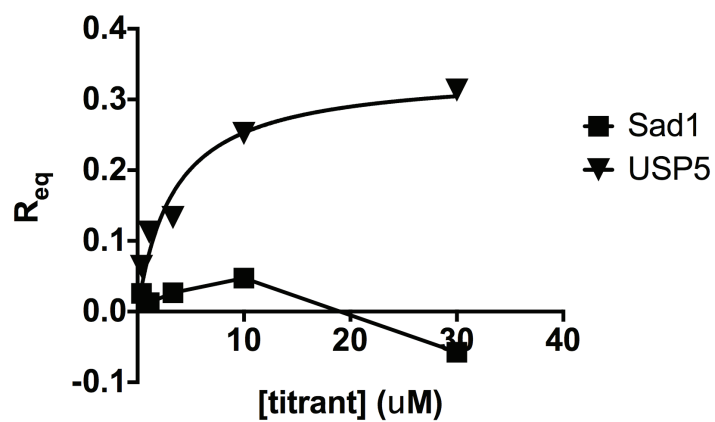
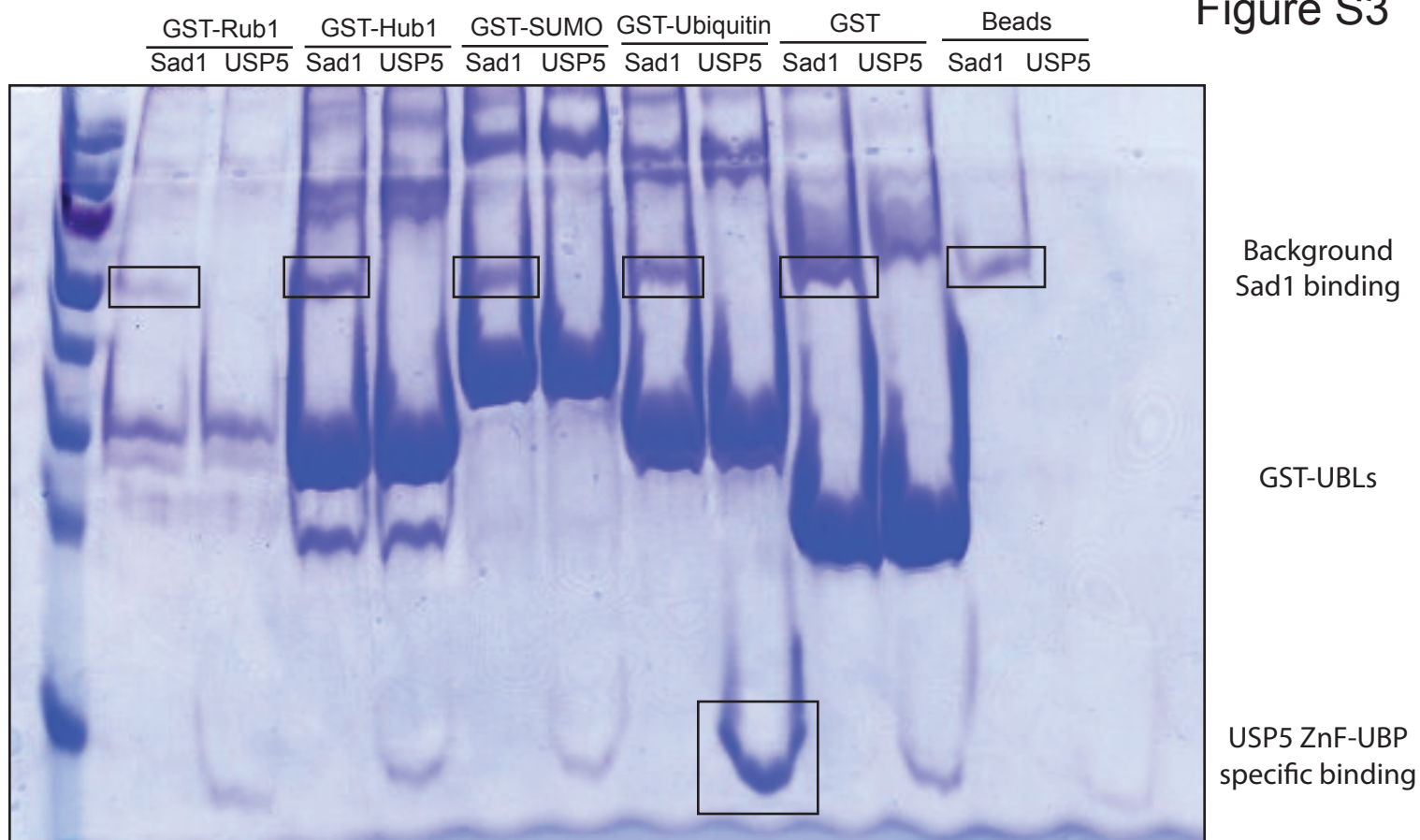
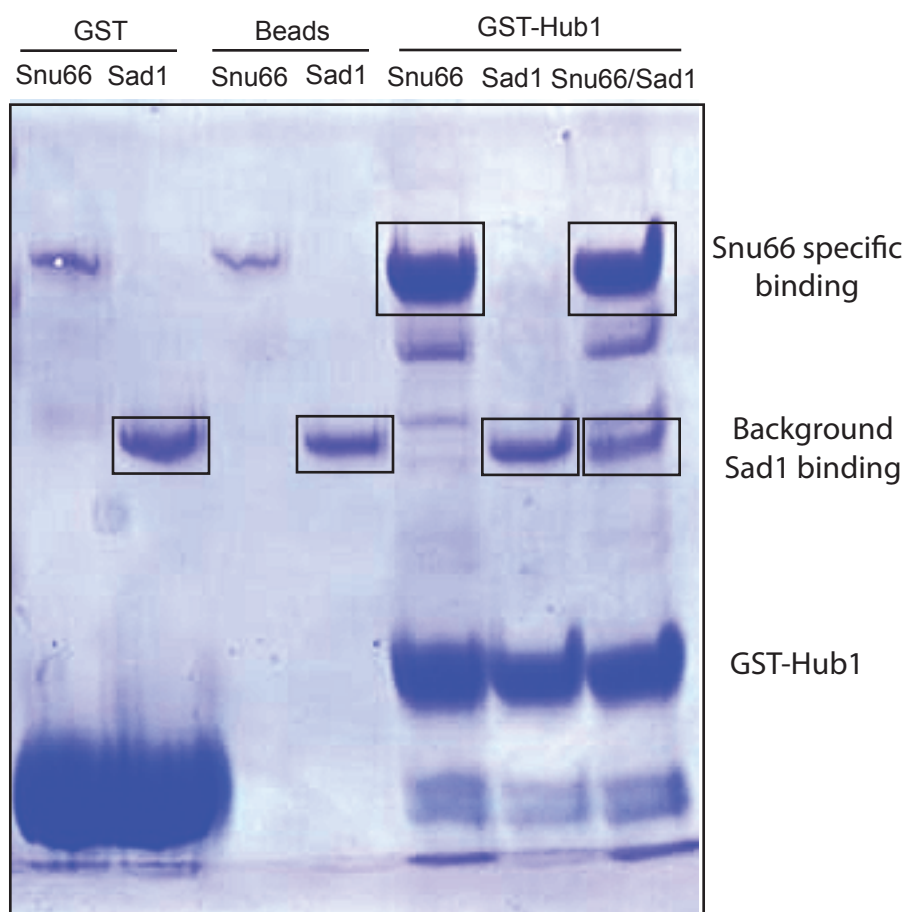


Figure S3

A



B



A

Sad1 Molprobity

Summary statistics

All-Atom Contacts	Clashscore, all atoms:	1.91	100 th percentile* (N=778, 1.870Å ± 0.25Å)	
	Clashscore is the number of serious steric overlaps (> 0.4 Å) per 1000 atoms.			
Protein Geometry	Poor rotamers	3	0.81%	Goal: <1%
	Ramachandran outliers	0	0.00%	Goal: <0.05%
	Ramachandran favored	378	97.67%	Goal: >98%
	MolProbity score*	1.03	100 th percentile* (N=11964, 1.870Å ± 0.25Å)	
	Cβ deviations >0.25Å	0	0.00%	Goal: 0
	Bad backbone bonds:	0 / 1576	0.00%	Goal: 0%
	Bad backbone angles:	0 / 1963	0.00%	Goal: <0.1%

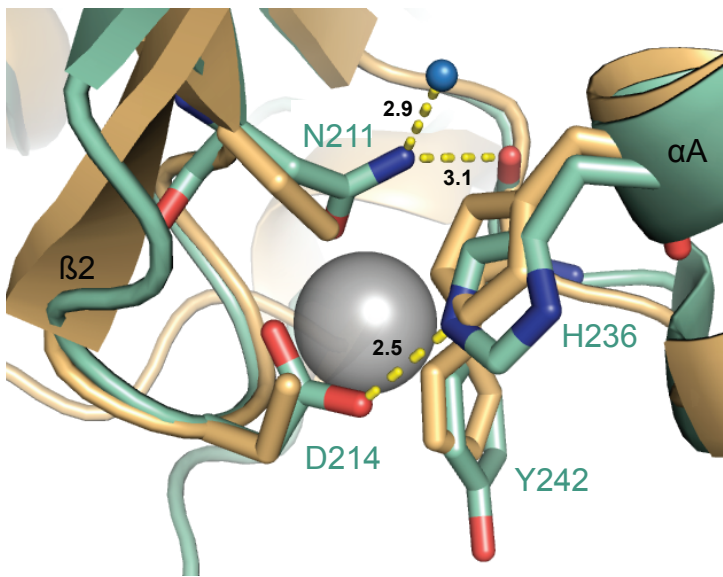
In the two column results, the left column gives the raw count, right column gives the percentage.

* 100th percentile is the best among structures of comparable resolution; 0th percentile is the worst. For clashscore the comparative set of structures was selected in 2004, for MolProbity score in 2006.

[†] MolProbity score combines the clashscore, rotamer, and Ramachandran evaluations into a single score, normalized to be on the same scale as X-ray resolution.

A

Sad1 ZnF-UBP
USP5 ZnF-UBP



B

USP5 homologues

	N211	D214		H236	Y242
<i>H. sapiens</i>	*	*		*	*
<i>M. musculus</i>					
<i>A. thaliana</i>					
<i>C. elegans</i>					
<i>D. melanogaster</i>					
<i>R. norvegicus</i>					
<i>S. pombe</i>					
<i>S. cerevisiae</i>					
Sad1					

NLWLNLTDG	SIL	CGRRYFDG	----	SGGNNHAVEHYRE	-TGYP
NLWLNLTDG	SIL	CGRRYFDG	----	SGGNNHAVEHYRE	-TGYP
NLWLNLTDG	MIL	CGRKNWDG	----	TGGNNHAVEHYKE	-TAYP
NLWLNLTDG	AVRCGR	SQFLSDG	KKTNGNGHM	QDYFDS	-TRFP
NLWLNLTDG	SIM	CGRKFFDG	----	SGGNDHAVEHYRV	-TGFP
NLWLNLTDG	SIL	CGRRYFDG	----	SGGNNHAVEHYRE	-TGYP
NLWMCLTCG	ALSCGR	KQYGG	----	GGGNGHALSHYDD	-TGHP
NLWLC	LHCGNIG	CGREQI	-G----	IDGHSHALDHYRSNNHP	
NVYCCLV	CGHY	YQGR	-HEKS	PAFIHSID	ENHH