



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:42 PM UTC

PDB ID : 2X6H / pdb_00002x6h
Title : THE CRYSTAL STRUCTURE OF THE DROSOPHILA CLASS III PI3-KINASE VPS34
Authors : Miller, S.; Tavshanjian, B.; Oleksy, A.; Perisic, O.; Houseman, B.T.; Shokat, K.M.; Williams, R.L.
Deposited on : 2010-02-17
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

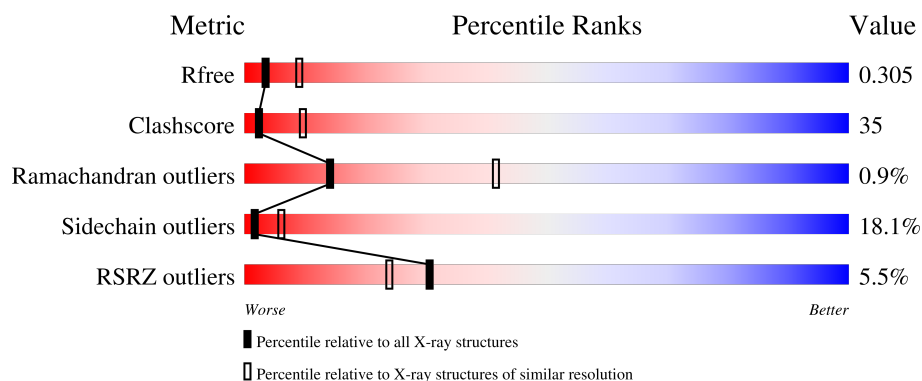
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	696	5% 35% 30% 11% • 22%
1	B	696	4% 36% 33% 9% 22%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8916 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called PHOSPHOTIDYLINOSITOL 3 KINASE 59F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	543	Total	C	N	O	S	0	0	0
			4445	2874	755	789	27			
1	B	543	Total	C	N	O	S	0	0	0
			4441	2870	759	785	27			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	254	GLY	-	expression tag	UNP Q9W1M7
A	255	SER	-	expression tag	UNP Q9W1M7
A	256	HIS	-	expression tag	UNP Q9W1M7
A	257	MET	-	expression tag	UNP Q9W1M7
A	455	ALA	GLY	engineered mutation	UNP Q9W1M7
B	254	GLY	-	expression tag	UNP Q9W1M7
B	255	SER	-	expression tag	UNP Q9W1M7
B	256	HIS	-	expression tag	UNP Q9W1M7
B	257	MET	-	expression tag	UNP Q9W1M7
B	455	ALA	GLY	engineered mutation	UNP Q9W1M7

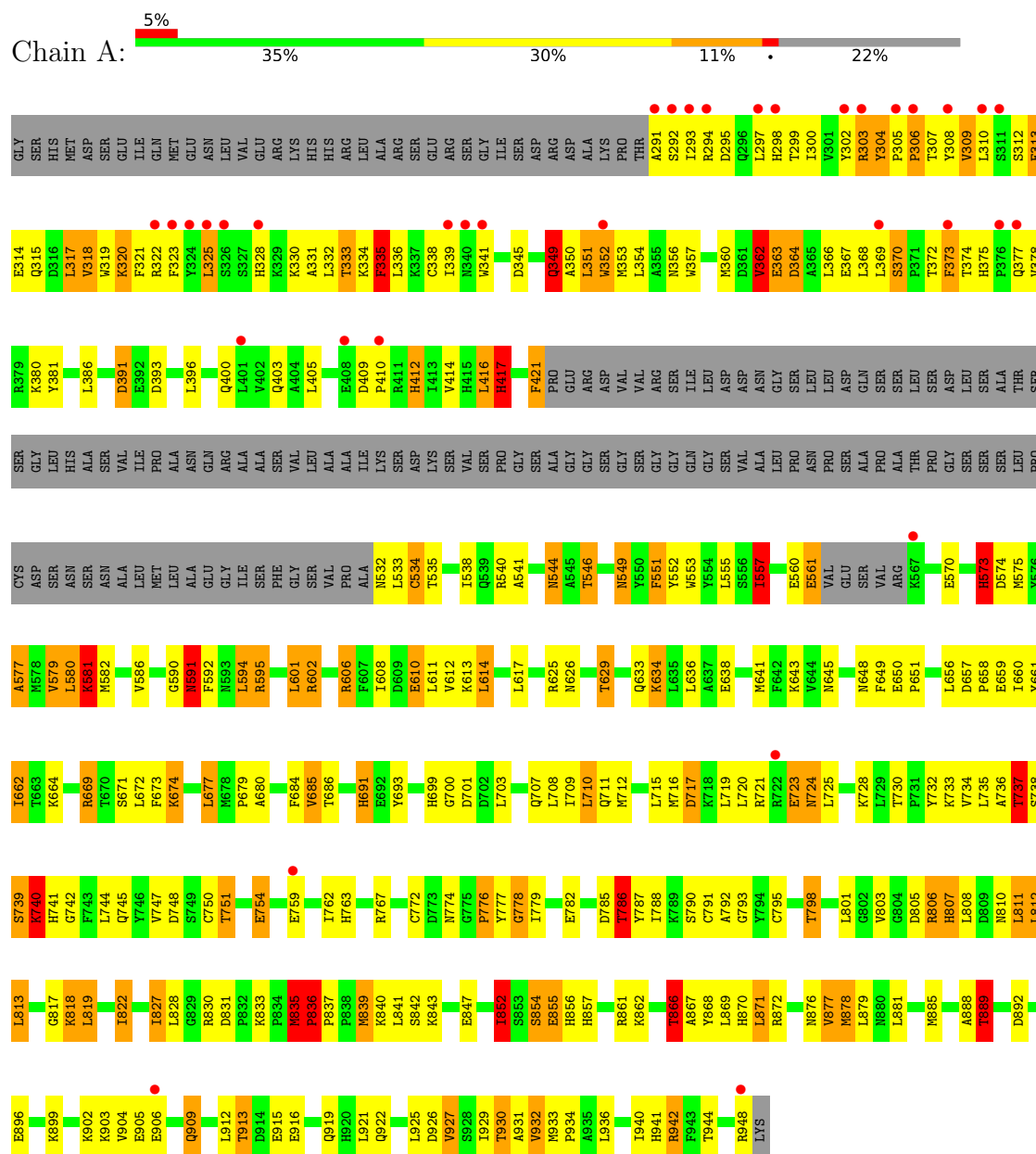
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	11	Total	O	0	0
			11	11		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PHOSPHOTIDYLINOSITOL 3 KINASE 59F



• Molecule 1: PHOSPHOTIDYLINOSITOL 3 KINASE 59F



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	110.60Å 154.79Å 243.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	72.36 – 2.90 72.36 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.7 (72.36-2.90) 95.6 (72.36-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.240 , 0.278 (Not available) , 0.305	Depositor DCC
R_{free} test set	2228 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	76.2	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 73.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8916	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.74	1/4553 (0.0%)	1.45	83/6164 (1.3%)
1	B	0.59	0/4548	1.04	8/6156 (0.1%)
All	All	0.67	1/9101 (0.0%)	1.26	91/12320 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	889	THR	CA-CB	5.15	1.61	1.53

All (91) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	592	PHE	N-CA-C	-10.86	100.16	113.41
1	A	835	MET	N-CA-C	9.11	121.25	110.31
1	A	380	LYS	N-CA-C	-8.98	101.50	111.28
1	A	318	VAL	N-CA-C	-8.73	101.72	110.62
1	B	343	LEU	N-CA-C	8.27	120.21	111.03
1	A	557	ILE	N-CA-C	-8.08	101.00	111.09
1	A	535	THR	N-CA-C	-7.92	102.63	111.82
1	A	684	PHE	CA-C-N	-7.62	111.75	122.34
1	A	684	PHE	C-N-CA	-7.62	111.75	122.34
1	A	577	ALA	N-CA-C	-7.42	103.19	111.28
1	A	735	LEU	CA-C-N	-7.39	111.64	122.65
1	A	735	LEU	C-N-CA	-7.39	111.64	122.65
1	A	944	THR	N-CA-C	-7.20	103.44	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304	TYR	CA-C-N	-7.17	112.99	120.38
1	B	304	TYR	C-N-CA	-7.17	112.99	120.38
1	A	818	LYS	N-CA-C	-6.95	100.20	110.48
1	A	836	PRO	N-CA-C	-6.94	102.23	110.70
1	A	369	LEU	N-CA-C	-6.89	104.34	112.89
1	A	866	THR	N-CA-C	-6.80	103.66	112.23
1	A	798	THR	N-CA-C	-6.78	103.68	112.23
1	A	793	GLY	N-CA-C	-6.77	104.10	112.77
1	A	551	PHE	N-CA-C	-6.76	103.91	111.28
1	A	915	GLU	N-CA-C	-6.76	103.99	111.36
1	A	403	GLN	N-CA-C	-6.58	104.19	111.82
1	A	867	ALA	N-CA-C	-6.55	104.22	111.36
1	A	852	ILE	N-CA-C	6.49	117.18	110.36
1	A	617	LEU	N-CA-C	-6.33	104.46	111.36
1	A	693	TYR	CA-C-N	-6.31	112.72	122.87
1	A	693	TYR	C-N-CA	-6.31	112.72	122.87
1	A	717	ASP	N-CA-C	-6.29	104.50	111.36
1	A	309	VAL	N-CA-C	-6.22	104.61	113.07
1	A	871	LEU	N-CA-C	-6.17	104.63	111.36
1	B	344	GLU	N-CA-C	6.17	123.94	110.80
1	A	677	LEU	N-CA-C	-6.03	101.36	110.28
1	A	787	TYR	N-CA-C	-6.02	104.72	111.28
1	A	352	TRP	N-CA-C	-5.98	104.84	111.36
1	A	785	ASP	N-CA-C	-5.89	104.94	111.36
1	A	573	HIS	N-CA-C	-5.88	104.77	111.07
1	A	921	LEU	N-CA-C	-5.86	104.98	111.36
1	A	792	ALA	N-CA-C	-5.85	104.49	111.69
1	A	786	THR	N-CA-CB	5.84	119.30	110.30
1	A	650	GLU	CA-C-N	5.79	126.12	119.93
1	A	650	GLU	C-N-CA	5.79	126.12	119.93
1	A	740	LYS	CA-C-N	-5.75	113.66	122.62
1	A	740	LYS	C-N-CA	-5.75	113.66	122.62
1	A	333	THR	N-CA-C	-5.71	105.05	111.28
1	A	906	GLU	N-CA-C	-5.70	105.14	111.36
1	A	416	LEU	N-CA-C	-5.68	105.17	111.36
1	A	878	MET	N-CA-C	-5.68	105.17	111.36
1	A	774	ASN	N-CA-C	5.67	117.54	111.36
1	A	927	VAL	N-CA-C	-5.67	104.99	110.72
1	A	579	VAL	N-CA-C	-5.66	105.00	110.72
1	A	610	GLU	N-CA-C	-5.66	104.73	111.69
1	A	581	LYS	N-CA-C	-5.65	105.27	111.82
1	A	325	LEU	N-CA-C	5.64	117.57	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	345	ASP	N-CA-C	-5.59	105.34	111.82
1	A	594	LEU	N-CA-C	-5.57	104.71	112.45
1	A	417	HIS	N-CA-C	-5.56	105.22	111.28
1	A	595	ARG	N-CA-C	-5.53	105.33	111.36
1	A	707	GLN	N-CA-C	-5.53	105.33	111.36
1	A	335	PHE	N-CA-C	-5.51	105.27	111.28
1	A	739	SER	N-CA-C	-5.50	102.25	110.23
1	B	888	ALA	N-CA-C	5.49	118.19	111.82
1	A	351	LEU	N-CA-C	-5.47	105.47	111.82
1	A	801	LEU	N-CA-C	-5.41	106.73	113.38
1	A	833	LYS	CA-C-N	5.39	125.06	119.56
1	A	833	LYS	C-N-CA	5.39	125.06	119.56
1	B	678	MET	N-CA-C	5.38	118.38	108.69
1	A	701	ASP	N-CA-C	5.38	116.81	109.18
1	A	904	VAL	N-CA-C	-5.31	105.36	110.72
1	A	364	ASP	N-CA-C	-5.30	105.08	112.45
1	A	941	HIS	N-CA-C	-5.24	105.57	111.28
1	A	927	VAL	CB-CA-C	-5.24	104.99	112.22
1	A	836	PRO	CA-C-N	5.24	139.57	127.00
1	A	836	PRO	C-N-CA	5.24	139.57	127.00
1	A	778	GLY	N-CA-C	-5.22	107.78	115.30
1	B	326	SER	N-CA-C	-5.21	106.12	113.20
1	A	817	GLY	CA-C-N	-5.20	112.76	121.39
1	A	817	GLY	C-N-CA	-5.20	112.76	121.39
1	A	538	ILE	N-CA-C	-5.19	105.48	110.72
1	A	320	LYS	N-CA-C	-5.16	105.83	111.82
1	A	836	PRO	C-N-CD	-5.16	109.25	120.60
1	B	305	PRO	N-CA-C	5.14	116.97	110.70
1	A	720	LEU	N-CA-C	-5.11	105.79	111.36
1	A	541	ALA	N-CA-C	-5.07	105.88	111.71
1	A	754	GLU	N-CA-C	-5.07	105.83	111.36
1	A	601	LEU	N-CA-C	-5.07	105.88	111.71
1	A	680	ALA	N-CA-C	-5.05	103.00	110.48
1	A	349	GLN	N-CA-C	-5.05	105.87	112.23
1	A	330	LYS	N-CA-C	-5.00	106.02	111.82
1	A	868	TYR	N-CA-C	-5.00	105.91	111.36

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	590	GLY	Mainchain

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Mol	Chain	Res	Type	Group
1	A	591	ASN	Mainchain
1	A	737	THR	Mainchain
1	A	835	MET	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4445	0	4478	284	1
1	B	4441	0	4485	353	0
2	A	19	0	0	4	0
2	B	11	0	0	0	0
All	All	8916	0	8963	630	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 35.

All (630) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:602:ARG:HD2	1:A:606:ARG:NH1	1.38	1.34
1:A:335:PHE:CE2	1:A:339:ILE:HD11	1.70	1.27
1:A:322:ARG:NH2	1:A:338:CYS:SG	2.11	1.22
1:A:335:PHE:HD1	1:A:357:TRP:CZ3	1.58	1.21
1:A:335:PHE:CD1	1:A:357:TRP:CZ3	2.30	1.19
1:B:669:ARG:NH1	1:B:683:THR:OG1	1.84	1.10
1:B:738:SER:HB3	1:B:741:HIS:CE1	1.86	1.10
1:B:315:GLN:HB2	1:B:338:CYS:HB3	1.31	1.10
1:A:335:PHE:CD1	1:A:357:TRP:HZ3	1.65	1.09
1:B:354:LEU:HD21	1:B:381:TYR:CD1	1.88	1.07
1:A:409:ASP:HB2	1:A:412:HIS:CD2	1.90	1.06
1:A:319:TRP:HA	1:A:322:ARG:HD2	1.39	1.04
1:A:669:ARG:HD3	2:A:2009:HOH:O	1.57	1.04
1:B:300:ILE:HA	1:B:304:TYR:CE2	1.92	1.04
1:A:335:PHE:CD2	1:A:339:ILE:HD11	1.94	1.03
1:A:306:PRO:HD2	1:A:308:TYR:CE2	1.95	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:602:ARG:CD	1:A:606:ARG:HH12	1.75	0.99
1:A:409:ASP:HB2	1:A:412:HIS:NE2	1.76	0.99
1:A:291:ALA:HB3	1:A:294:ARG:HB2	1.45	0.99
1:A:335:PHE:HE2	1:A:339:ILE:HD11	1.20	0.98
1:A:577:ALA:O	1:A:581:LYS:HD2	1.64	0.97
1:A:602:ARG:CD	1:A:606:ARG:NH1	2.30	0.94
1:B:739:SER:O	1:B:740:LYS:HG2	1.67	0.94
1:B:639:GLN:HE21	1:B:645:ASN:HD21	1.11	0.94
1:B:738:SER:HB3	1:B:741:HIS:HE1	1.30	0.92
1:B:837:PRO:O	1:B:840:LYS:HE2	1.70	0.92
1:B:300:ILE:HA	1:B:304:TYR:HE2	1.28	0.91
1:B:375:HIS:CD2	1:B:377:GLN:H	1.89	0.91
1:A:734:VAL:HG22	1:A:744:LEU:HD23	1.52	0.91
1:B:297:LEU:O	1:B:300:ILE:HG22	1.71	0.91
1:B:300:ILE:O	1:B:304:TYR:HD2	1.53	0.91
1:B:322:ARG:HH22	1:B:334:LYS:HB2	1.35	0.91
1:A:806:ARG:NH2	1:A:822:ILE:O	2.03	0.90
1:B:639:GLN:NE2	1:B:645:ASN:HD21	1.67	0.90
1:A:306:PRO:HD2	1:A:308:TYR:CD2	2.06	0.90
1:A:319:TRP:HA	1:A:322:ARG:CD	2.03	0.88
1:B:375:HIS:HD2	1:B:377:GLN:H	1.15	0.88
1:B:300:ILE:O	1:B:304:TYR:CD2	2.26	0.87
1:B:782:GLU:O	1:B:786:THR:HG22	1.73	0.87
1:A:608:ILE:O	1:A:612:VAL:HG23	1.73	0.87
1:B:836:PRO:HB2	1:B:837:PRO:HA	1.57	0.86
1:A:323:PHE:HA	1:A:353:MET:HE1	1.57	0.86
1:B:417:HIS:ND1	1:B:578:MET:HG3	1.90	0.85
1:A:335:PHE:HE2	1:A:339:ILE:CD1	1.88	0.85
1:B:633:GLN:NE2	1:B:668:MET:SD	2.50	0.85
1:A:367:GLU:O	1:A:370:SER:HB3	1.75	0.84
1:B:339:ILE:HD13	1:B:350:ALA:CB	2.08	0.84
1:A:625:ARG:O	1:A:629:THR:HG22	1.78	0.84
1:B:339:ILE:HD13	1:B:350:ALA:HB1	1.60	0.84
1:B:639:GLN:HE21	1:B:645:ASN:ND2	1.75	0.84
1:B:315:GLN:CB	1:B:338:CYS:HB3	2.06	0.83
1:B:751:THR:HG22	1:B:754:GLU:H	1.44	0.82
1:A:320:LYS:HG3	1:A:321:PHE:HD2	1.42	0.82
1:A:335:PHE:HD1	1:A:357:TRP:CH2	1.97	0.82
1:B:776:PRO:HB2	1:B:777:TYR:HD1	1.45	0.81
1:B:710:LEU:HD22	1:B:732:TYR:O	1.80	0.81
1:A:339:ILE:HG21	1:A:350:ALA:CB	2.11	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:LYS:HD3	1:B:361:ASP:OD1	1.80	0.81
1:B:812:LEU:HD22	1:B:812:LEU:N	1.96	0.81
1:B:311:SER:HA	1:B:314:GLU:CD	2.07	0.80
1:B:648:ASN:HA	1:B:664:LYS:HB3	1.63	0.80
1:A:674:LYS:H	1:A:674:LYS:HD3	1.45	0.79
1:A:335:PHE:CE2	1:A:339:ILE:CD1	2.58	0.79
1:A:602:ARG:HD2	1:A:606:ARG:HH12	0.80	0.79
1:A:323:PHE:HZ	1:A:349:GLN:HE22	1.32	0.78
1:B:763:HIS:CE1	1:B:777:TYR:CD2	2.71	0.78
1:B:667:PRO:O	1:B:670:THR:HG23	1.84	0.78
1:B:830:ARG:HH22	1:B:903:LYS:HZ2	1.31	0.78
1:B:888:ALA:O	1:B:889:THR:HG22	1.82	0.78
1:B:311:SER:O	1:B:315:GLN:N	2.16	0.77
1:B:315:GLN:HB2	1:B:338:CYS:CB	2.10	0.77
1:B:763:HIS:HE1	1:B:777:TYR:CD2	2.01	0.77
1:A:798:THR:HG23	1:A:803:VAL:HB	1.66	0.77
1:A:922:GLN:HE22	1:B:919:GLN:HB3	1.50	0.77
1:B:639:GLN:NE2	1:B:647:THR:OG1	2.18	0.77
1:B:824:PHE:HB3	1:B:827:ILE:HD11	1.65	0.77
1:A:306:PRO:CD	1:A:308:TYR:CE2	2.67	0.77
1:B:360:MET:HG2	1:B:364:ASP:HB2	1.65	0.77
1:B:716:MET:HG2	1:B:878:MET:HE2	1.67	0.77
1:A:827:ILE:HG23	1:A:828:LEU:HG	1.65	0.76
1:A:307:THR:OG1	1:A:905:GLU:OE2	2.03	0.76
1:A:717:ASP:OD2	1:A:721:ARG:NH1	2.19	0.76
1:B:398:LEU:HA	1:B:401:LEU:HB2	1.68	0.76
1:B:577:ALA:O	1:B:581:LYS:HD3	1.86	0.76
1:A:710:LEU:HD23	1:A:732:TYR:O	1.85	0.75
1:B:311:SER:HA	1:B:314:GLU:HB2	1.68	0.75
1:B:798:THR:HG23	1:B:803:VAL:HB	1.68	0.75
1:B:876:ASN:ND2	1:B:876:ASN:H	1.82	0.75
1:A:320:LYS:HG3	1:A:321:PHE:CD2	2.21	0.75
1:A:299:THR:HA	1:A:302:TYR:HD2	1.52	0.74
1:A:552:TYR:C	1:A:552:TYR:CD2	2.65	0.74
1:B:836:PRO:HB2	1:B:837:PRO:CA	2.17	0.74
1:A:412:HIS:CD2	1:A:412:HIS:H	2.05	0.74
1:A:677:LEU:O	1:A:679:PRO:HD3	1.87	0.74
1:B:667:PRO:O	1:B:670:THR:CG2	2.35	0.74
1:A:339:ILE:HG21	1:A:350:ALA:HB2	1.68	0.74
1:A:608:ILE:HG21	1:A:737:THR:HG21	1.68	0.73
1:A:861:ARG:NH2	1:A:926:ASP:OD2	2.20	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:LEU:HB3	1:A:581:LYS:HE3	1.70	0.73
1:B:363:GLU:HG2	1:B:877:VAL:HG22	1.70	0.73
1:A:294:ARG:NH2	1:A:298:HIS:NE2	2.36	0.73
1:B:306:PRO:HG2	1:B:879:LEU:HD12	1.70	0.73
1:B:751:THR:CG2	1:B:754:GLU:H	2.02	0.73
1:B:329:LYS:HG2	1:B:359:PRO:O	1.87	0.73
1:B:360:MET:HE2	1:B:381:TYR:HE2	1.54	0.72
1:A:872:ARG:NH1	1:A:909:GLN:O	2.22	0.72
1:B:302:TYR:CE1	1:B:331:ALA:HB2	2.24	0.72
1:A:295:ASP:O	1:A:299:THR:HG23	1.90	0.71
1:A:306:PRO:HD2	1:A:308:TYR:HE2	1.51	0.71
1:A:306:PRO:CD	1:A:308:TYR:HE2	2.04	0.71
1:A:318:VAL:O	1:A:322:ARG:HG3	1.91	0.71
1:A:835:MET:HB3	1:A:836:PRO:O	1.91	0.71
1:A:930:THR:HG22	1:A:936:LEU:HB3	1.73	0.71
1:A:573:HIS:C	1:A:573:HIS:CD2	2.69	0.71
1:B:354:LEU:HD21	1:B:381:TYR:CE1	2.26	0.70
1:A:931:ALA:O	1:B:944:THR:OG1	2.09	0.70
1:B:416:LEU:HD22	1:B:535:THR:HG22	1.73	0.70
1:B:577:ALA:O	1:B:581:LYS:CD	2.39	0.70
1:A:669:ARG:CD	2:A:2009:HOH:O	2.28	0.70
1:A:926:ASP:O	1:A:930:THR:OG1	2.09	0.70
1:B:830:ARG:NH2	1:B:903:LYS:HZ2	1.89	0.70
1:A:948:ARG:HD2	1:A:948:ARG:C	2.16	0.70
1:B:302:TYR:HE1	1:B:331:ALA:HB2	1.55	0.70
1:B:722:ARG:HH11	1:B:722:ARG:HA	1.55	0.70
1:A:335:PHE:HD2	1:A:335:PHE:C	1.99	0.69
1:A:709:ILE:HG13	1:A:827:ILE:HD11	1.75	0.69
1:B:354:LEU:C	1:B:354:LEU:HD23	2.17	0.69
1:A:314:GLU:HA	1:A:317:LEU:HG	1.75	0.69
1:B:722:ARG:HH11	1:B:722:ARG:CA	2.04	0.69
1:B:812:LEU:HD22	1:B:812:LEU:H	1.58	0.69
1:A:335:PHE:CD2	1:A:335:PHE:C	2.71	0.69
1:A:314:GLU:O	1:A:318:VAL:HG23	1.93	0.69
1:A:335:PHE:CE1	1:A:357:TRP:HZ3	2.09	0.69
1:B:634:LYS:HA	1:B:634:LYS:HE2	1.74	0.69
1:B:664:LYS:HG3	1:B:685:VAL:CG2	2.24	0.68
1:A:866:THR:O	1:A:870:HIS:HD2	1.76	0.68
1:A:291:ALA:HB3	1:A:294:ARG:CB	2.21	0.68
1:B:703:LEU:HB2	1:B:736:ALA:HB2	1.76	0.68
1:B:391:ASP:OD2	1:B:540:ARG:NH2	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:629:THR:HB	1:A:672:LEU:HG	1.76	0.67
1:A:677:LEU:HD22	1:A:700:GLY:HA3	1.76	0.67
1:B:354:LEU:HD21	1:B:381:TYR:HD1	1.58	0.67
1:B:400:GLN:HE22	1:B:711:GLN:HE22	1.42	0.67
1:B:416:LEU:HB3	1:B:582:MET:HE1	1.75	0.67
1:B:331:ALA:HB1	1:B:357:TRP:HZ2	1.60	0.67
1:A:909:GLN:HA	1:A:909:GLN:OE1	1.93	0.67
1:A:806:ARG:HD2	1:A:806:ARG:N	2.09	0.67
1:A:913:THR:HG23	1:A:916:GLU:HB2	1.76	0.67
1:B:843:LYS:HA	1:B:932:VAL:CG1	2.25	0.67
1:B:322:ARG:HH22	1:B:334:LYS:CB	2.08	0.67
1:B:370:SER:OG	1:B:372:THR:HG23	1.94	0.66
1:A:352:TRP:O	1:A:356:ASN:ND2	2.28	0.66
1:B:360:MET:CG	1:B:364:ASP:HB2	2.26	0.66
1:B:664:LYS:CG	1:B:685:VAL:CG2	2.74	0.66
1:B:776:PRO:HD2	1:B:779:ILE:O	1.94	0.66
1:A:373:PHE:N	1:A:373:PHE:HD2	1.94	0.66
1:B:876:ASN:H	1:B:876:ASN:HD22	1.43	0.66
1:A:777:TYR:HB2	1:A:779:ILE:HD12	1.78	0.66
1:B:716:MET:HG2	1:B:878:MET:CE	2.26	0.65
1:A:300:ILE:O	1:A:304:TYR:HB2	1.96	0.65
1:B:863:GLN:O	1:B:866:THR:HG22	1.96	0.65
1:B:764:ASN:OD1	1:B:767:ARG:NH1	2.29	0.65
1:B:823:ASP:OD1	1:B:823:ASP:C	2.40	0.65
1:A:409:ASP:CB	1:A:412:HIS:CD2	2.73	0.65
1:B:654:PHE:HD1	1:B:655:PRO:HD2	1.61	0.65
1:A:737:THR:HG22	1:A:738:SER:H	1.61	0.65
1:B:716:MET:CG	1:B:878:MET:HE2	2.27	0.65
1:A:306:PRO:CD	1:A:308:TYR:CD2	2.80	0.64
1:B:553:TRP:O	1:B:557:ILE:HG13	1.96	0.64
1:A:360:MET:HE2	1:A:364:ASP:HB3	1.80	0.64
1:B:698:LYS:O	1:B:741:HIS:HA	1.97	0.64
1:B:751:THR:HG22	1:B:754:GLU:CB	2.27	0.64
1:B:830:ARG:HH22	1:B:903:LYS:NZ	1.95	0.64
1:A:737:THR:HG22	1:A:738:SER:N	2.13	0.64
1:B:733:LYS:H	1:B:745:GLN:HE21	1.45	0.64
1:A:405:LEU:HD23	1:A:533:LEU:CD2	2.26	0.64
1:B:751:THR:HG22	1:B:754:GLU:HB2	1.80	0.64
1:B:830:ARG:HH12	1:B:903:LYS:HZ2	1.46	0.63
1:B:836:PRO:CB	1:B:837:PRO:CA	2.76	0.63
1:B:375:HIS:HD2	1:B:377:GLN:N	1.93	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:PHE:CE1	1:A:349:GLN:OE1	2.50	0.63
1:A:625:ARG:O	1:A:629:THR:CG2	2.46	0.63
1:B:669:ARG:HH11	1:B:683:THR:HG1	1.43	0.63
1:A:656:LEU:HD22	1:A:733:LYS:HB3	1.81	0.63
1:B:314:GLU:O	1:B:317:LEU:HG	1.99	0.63
1:B:836:PRO:HB3	1:B:837:PRO:C	2.24	0.62
1:B:417:HIS:ND1	1:B:578:MET:CG	2.61	0.62
1:A:412:HIS:H	1:A:412:HIS:HD2	1.45	0.62
1:B:318:VAL:O	1:B:322:ARG:HG3	2.00	0.62
1:A:400:GLN:HG2	1:A:881:LEU:HD22	1.82	0.62
1:A:561:GLU:CD	1:A:561:GLU:H	2.07	0.62
1:A:575:MET:O	1:A:579:VAL:HG23	1.99	0.61
1:A:373:PHE:N	1:A:373:PHE:CD2	2.66	0.61
1:B:568:GLN:OE1	1:B:568:GLN:HA	1.99	0.61
1:A:417:HIS:HD1	1:A:417:HIS:C	2.08	0.61
1:A:710:LEU:CD2	1:A:732:TYR:O	2.48	0.61
1:B:777:TYR:HB2	1:B:779:ILE:HD12	1.83	0.61
1:A:942:ARG:HD3	1:B:835:MET:HG3	1.82	0.61
1:B:746:TYR:C	1:B:746:TYR:CD2	2.79	0.61
1:A:553:TRP:O	1:A:557:ILE:HG13	2.01	0.60
1:A:888:ALA:O	1:A:889:THR:HG23	2.01	0.60
1:B:810:ASN:O	1:B:822:ILE:HG22	2.02	0.60
1:A:674:LYS:HD3	1:A:674:LYS:N	2.16	0.60
1:B:557:ILE:O	1:B:560:GLU:HG2	2.02	0.60
1:A:751:THR:HG23	1:A:754:GLU:CG	2.31	0.60
1:B:363:GLU:HG3	1:B:877:VAL:HG13	1.83	0.60
1:A:294:ARG:HH22	1:A:317:LEU:HB2	1.66	0.60
1:B:898:ASP:OD1	1:B:898:ASP:N	2.34	0.60
1:B:319:TRP:NE1	1:B:323:PHE:CZ	2.69	0.60
1:B:664:LYS:CG	1:B:685:VAL:HG21	2.32	0.59
1:B:780:SER:HB3	1:B:782:GLU:OE2	2.02	0.59
1:A:313:GLU:CA	1:A:313:GLU:OE1	2.49	0.59
1:A:629:THR:O	1:A:633:GLN:HG3	2.02	0.59
1:B:812:LEU:N	1:B:812:LEU:CD2	2.65	0.59
1:A:322:ARG:HH22	1:A:338:CYS:HG	1.43	0.59
1:A:674:LYS:H	1:A:674:LYS:CD	2.09	0.59
1:B:645:ASN:OD1	1:B:645:ASN:C	2.45	0.59
1:B:830:ARG:NH1	1:B:903:LYS:HZ2	2.00	0.59
1:A:322:ARG:HH12	1:A:338:CYS:HB2	1.68	0.59
1:A:813:LEU:HD12	1:A:818:LYS:O	2.02	0.59
1:B:348:THR:HA	1:B:351:LEU:HD22	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:SER:HB2	1:B:371:PRO:CD	2.33	0.59
1:B:636:LEU:CD1	1:B:670:THR:HG21	2.33	0.59
1:B:677:LEU:O	1:B:679:PRO:HD3	2.02	0.59
1:B:768:LYS:HB3	1:B:769:HIS:HD2	1.68	0.58
1:A:533:LEU:C	1:A:533:LEU:HD23	2.28	0.58
1:A:827:ILE:HG22	1:A:892:ASP:CB	2.32	0.58
1:A:739:SER:O	1:A:740:LYS:HG2	2.03	0.58
1:B:325:LEU:C	1:B:327:SER:H	2.11	0.58
1:B:360:MET:HE2	1:B:381:TYR:CE2	2.38	0.58
1:B:320:LYS:HG3	1:B:321:PHE:N	2.19	0.58
1:B:541:ALA:O	1:B:548:ALA:HB2	2.03	0.58
1:A:772:CYS:O	1:A:778:GLY:HA2	2.03	0.58
1:B:836:PRO:CB	1:B:837:PRO:C	2.77	0.58
1:B:322:ARG:HD3	1:B:335:PHE:CD1	2.39	0.58
1:B:909:GLN:HG3	1:B:912:LEU:HG	1.85	0.58
1:B:639:GLN:NE2	1:B:645:ASN:ND2	2.43	0.58
1:B:568:GLN:OE1	1:B:571:ARG:NH1	2.37	0.58
1:B:558:GLU:HB3	1:B:576:TYR:CD1	2.38	0.57
1:B:803:VAL:HG12	1:B:806:ARG:HD3	1.85	0.57
1:B:732:TYR:HB2	1:B:745:GLN:HB3	1.86	0.57
1:B:727:LEU:HB2	1:B:729:LEU:HD21	1.85	0.57
1:A:391:ASP:OD2	1:A:391:ASP:N	2.36	0.57
1:B:722:ARG:HH11	1:B:722:ARG:CG	2.16	0.57
1:A:305:PRO:HA	1:A:308:TYR:CD2	2.40	0.57
1:A:306:PRO:HB3	1:A:876:ASN:OD1	2.05	0.57
1:B:322:ARG:NH2	1:B:334:LYS:HB2	2.13	0.57
1:B:699:HIS:CD2	1:B:699:HIS:H	2.22	0.57
1:B:664:LYS:HG2	1:B:685:VAL:CG2	2.35	0.56
1:A:854:SER:O	1:A:855:GLU:C	2.48	0.56
1:B:417:HIS:CE1	1:B:578:MET:HG3	2.39	0.56
1:B:696:ILE:HB	1:B:744:LEU:HB2	1.88	0.56
1:B:722:ARG:HH11	1:B:722:ARG:CB	2.18	0.56
1:A:811:LEU:C	1:A:812:LEU:HD23	2.31	0.56
1:B:300:ILE:CG1	1:B:304:TYR:CE2	2.89	0.56
1:B:715:LEU:O	1:B:719:LEU:HD23	2.06	0.56
1:B:885:MET:O	1:B:888:ALA:HB2	2.05	0.56
1:A:294:ARG:NH2	1:A:317:LEU:HB2	2.21	0.56
1:A:691:HIS:ND1	2:A:2008:HOH:O	1.98	0.56
1:A:582:MET:O	1:A:586:VAL:HG23	2.05	0.56
1:A:302:TYR:HD1	1:A:331:ALA:HA	1.71	0.56
1:A:751:THR:HG23	1:A:754:GLU:HG3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:805:ASP:O	1:A:810:ASN:ND2	2.38	0.56
1:B:751:THR:HG22	1:B:754:GLU:N	2.19	0.56
1:A:360:MET:HG2	1:A:364:ASP:HB2	1.87	0.56
1:B:300:ILE:CA	1:B:304:TYR:CE2	2.81	0.56
1:A:335:PHE:HD2	1:A:335:PHE:O	1.89	0.55
1:B:335:PHE:CD1	1:B:357:TRP:CZ3	2.93	0.55
1:B:830:ARG:HH12	1:B:903:LYS:NZ	2.03	0.55
1:A:313:GLU:OE1	1:A:313:GLU:HA	2.07	0.55
1:B:398:LEU:CD1	1:B:554:TYR:CE2	2.90	0.55
1:B:669:ARG:NH1	1:B:683:THR:CG2	2.70	0.55
1:A:318:VAL:O	1:A:322:ARG:N	2.39	0.55
1:A:339:ILE:C	1:A:341:TRP:H	2.11	0.55
1:A:354:LEU:HD21	1:A:381:TYR:CD1	2.41	0.55
1:A:919:GLN:HB3	1:B:922:GLN:HE22	1.71	0.55
1:B:400:GLN:HE22	1:B:711:GLN:NE2	2.05	0.55
1:A:669:ARG:NH2	2:A:2006:HOH:O	2.38	0.55
1:B:311:SER:HA	1:B:314:GLU:OE2	2.06	0.55
1:B:806:ARG:N	1:B:806:ARG:HD2	2.23	0.55
1:A:322:ARG:NH2	1:A:338:CYS:HG	2.02	0.54
1:A:354:LEU:HD23	1:A:354:LEU:O	2.07	0.54
1:B:837:PRO:O	1:B:840:LYS:CE	2.50	0.54
1:B:300:ILE:HG13	1:B:304:TYR:CE2	2.42	0.54
1:B:649:PHE:CE2	1:B:664:LYS:HA	2.42	0.54
1:A:657:ASP:HB3	1:A:660:ILE:HD12	1.88	0.54
1:A:803:VAL:HG12	1:A:806:ARG:HD3	1.90	0.54
1:A:854:SER:C	1:A:856:HIS:N	2.64	0.54
1:A:305:PRO:O	1:A:334:LYS:NZ	2.39	0.54
1:A:360:MET:HE1	1:A:368:LEU:HD21	1.89	0.54
1:A:942:ARG:HD3	1:B:835:MET:CG	2.38	0.54
1:B:738:SER:CB	1:B:741:HIS:HE1	2.13	0.54
1:A:909:GLN:HG3	1:A:912:LEU:HG	1.89	0.54
1:A:839:MET:HG2	1:A:925:LEU:HD23	1.90	0.54
1:A:843:LYS:HE2	1:A:847:GLU:OE2	2.06	0.54
1:A:879:LEU:HD11	1:A:905:GLU:HG3	1.90	0.54
1:A:319:TRP:O	1:A:322:ARG:HB2	2.08	0.54
1:A:677:LEU:CD2	1:A:700:GLY:HA3	2.38	0.54
1:A:737:THR:CG2	1:A:741:HIS:CD2	2.90	0.54
1:B:342:LYS:HB3	1:B:346:GLU:HG3	1.90	0.54
1:B:702:ASP:OD1	1:B:704:ARG:NH2	2.39	0.54
1:B:705:GLN:HG2	1:B:827:ILE:HD13	1.90	0.54
1:B:788:ILE:HG13	1:B:860:PHE:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:639:GLN:HG2	1:B:647:THR:HG21	1.89	0.53
1:B:730:THR:HG22	1:B:818:LYS:HD3	1.90	0.53
1:A:782:GLU:O	1:A:786:THR:HG23	2.08	0.53
1:B:306:PRO:HG2	1:B:879:LEU:CD1	2.39	0.53
1:B:313:GLU:HG2	1:B:314:GLU:HG3	1.90	0.53
1:B:669:ARG:HH11	1:B:683:THR:CB	2.19	0.53
1:B:322:ARG:HD3	1:B:335:PHE:HD1	1.73	0.53
1:B:369:LEU:HD11	1:B:386:LEU:HD11	1.90	0.53
1:B:631:LYS:HD2	1:B:635:LEU:CD2	2.37	0.53
1:A:614:LEU:HD21	1:A:636:LEU:HD23	1.91	0.53
1:A:776:PRO:HB2	1:A:777:TYR:HD1	1.73	0.53
1:B:306:PRO:CG	1:B:879:LEU:HD12	2.38	0.53
1:A:304:TYR:HB2	1:A:305:PRO:HD3	1.90	0.53
1:A:739:SER:O	1:A:740:LYS:CG	2.56	0.53
1:B:654:PHE:CD1	1:B:655:PRO:HD2	2.43	0.53
1:B:810:ASN:C	1:B:811:LEU:HD23	2.34	0.53
1:B:534:CYS:O	1:B:538:ILE:HG13	2.08	0.53
1:B:709:ILE:HD12	1:B:801:LEU:HD13	1.91	0.53
1:B:715:LEU:HD12	1:B:881:LEU:HD12	1.91	0.53
1:B:575:MET:C	1:B:575:MET:HE2	2.34	0.53
1:B:732:TYR:OH	1:B:822:ILE:O	2.25	0.53
1:A:417:HIS:C	1:A:417:HIS:ND1	2.67	0.52
1:A:580:LEU:HB3	1:A:581:LYS:CE	2.39	0.52
1:A:807:HIS:H	1:A:810:ASN:HB2	1.73	0.52
1:B:348:THR:HA	1:B:351:LEU:CD2	2.39	0.52
1:B:839:MET:HE1	1:B:921:LEU:HD11	1.91	0.52
1:A:782:GLU:O	1:A:786:THR:CG2	2.58	0.52
1:B:311:SER:OG	1:B:314:GLU:HB2	2.08	0.52
1:B:381:TYR:C	1:B:381:TYR:CD2	2.87	0.52
1:A:323:PHE:CA	1:A:353:MET:HE1	2.33	0.52
1:B:319:TRP:CD1	1:B:323:PHE:CE2	2.97	0.52
1:A:421:PHE:N	1:A:421:PHE:CD1	2.78	0.52
1:B:292:SER:O	1:B:295:ASP:HB3	2.10	0.52
1:A:405:LEU:HD23	1:A:533:LEU:HD21	1.90	0.52
1:B:719:LEU:HD23	1:B:719:LEU:N	2.25	0.52
1:B:806:ARG:CZ	1:B:810:ASN:HD22	2.23	0.52
1:A:410:PRO:O	1:A:414:VAL:HG23	2.10	0.52
1:A:560:GLU:OE2	1:A:739:SER:OG	2.27	0.52
1:A:649:PHE:HE1	1:A:662:ILE:HG13	1.73	0.52
1:B:410:PRO:O	1:B:414:VAL:HG22	2.10	0.52
1:B:551:PHE:O	1:B:555:LEU:HG	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:669:ARG:NH1	1:B:683:THR:CB	2.73	0.52
1:B:830:ARG:CZ	1:B:903:LYS:HZ2	2.23	0.52
1:B:319:TRP:NE1	1:B:323:PHE:CE2	2.78	0.51
1:B:335:PHE:C	1:B:335:PHE:CD2	2.88	0.51
1:A:712:MET:O	1:A:716:MET:HG3	2.10	0.51
1:B:311:SER:CA	1:B:314:GLU:HB2	2.38	0.51
1:B:575:MET:HE2	1:B:576:TYR:HA	1.91	0.51
1:B:577:ALA:O	1:B:581:LYS:HD2	2.11	0.51
1:B:823:ASP:OD1	1:B:824:PHE:N	2.43	0.51
1:B:839:MET:HG2	1:B:925:LEU:HD23	1.92	0.51
1:B:710:LEU:O	1:B:714:THR:OG1	2.26	0.51
1:B:737:THR:HG23	1:B:742:GLY:HA2	1.91	0.51
1:A:626:ASN:HA	1:A:629:THR:HG23	1.93	0.51
1:B:375:HIS:CD2	1:B:377:GLN:N	2.70	0.51
1:B:733:LYS:H	1:B:745:GLN:NE2	2.06	0.51
1:B:763:HIS:CD2	1:B:848:ALA:HA	2.45	0.51
1:B:827:ILE:CG2	1:B:828:LEU:HG	2.40	0.51
1:A:836:PRO:HB2	1:A:837:PRO:CA	2.41	0.51
1:B:664:LYS:HG3	1:B:685:VAL:HG21	1.89	0.51
1:A:354:LEU:HD23	1:A:354:LEU:C	2.36	0.51
1:A:405:LEU:CD2	1:A:533:LEU:CD2	2.88	0.51
1:B:708:LEU:HG	1:B:885:MET:HE2	1.92	0.51
1:B:822:ILE:HG23	1:B:823:ASP:N	2.26	0.51
1:A:363:GLU:HA	1:A:366:LEU:HD12	1.92	0.50
1:A:737:THR:HB	1:A:741:HIS:HD2	1.74	0.50
1:B:300:ILE:HG13	1:B:304:TYR:HE2	1.76	0.50
1:A:610:GLU:OE2	1:A:613:LYS:HD2	2.11	0.50
1:A:790:SER:CB	1:A:819:LEU:H	2.24	0.50
1:B:403:GLN:HE22	1:B:406:LYS:NZ	2.10	0.50
1:B:417:HIS:O	1:B:420:ILE:HG13	2.12	0.50
1:A:318:VAL:C	1:A:322:ARG:HG3	2.37	0.50
1:B:308:TYR:HD1	1:B:310:LEU:HD12	1.77	0.50
1:B:716:MET:HE2	1:B:878:MET:HE2	1.93	0.50
1:A:827:ILE:HG22	1:A:892:ASP:HB2	1.93	0.50
1:B:304:TYR:HB2	1:B:305:PRO:HD3	1.94	0.50
1:A:339:ILE:HD13	1:A:350:ALA:HB1	1.94	0.49
1:B:664:LYS:HG2	1:B:685:VAL:HG23	1.93	0.49
1:B:722:ARG:HH11	1:B:722:ARG:HG3	1.75	0.49
1:A:409:ASP:CB	1:A:412:HIS:NE2	2.64	0.49
1:A:579:VAL:HA	1:A:582:MET:HE3	1.94	0.49
1:B:321:PHE:O	1:B:325:LEU:HB2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:909:GLN:OE1	1:A:909:GLN:CA	2.60	0.49
1:B:810:ASN:HA	1:B:822:ILE:HG22	1.95	0.49
1:A:412:HIS:CD2	1:A:412:HIS:N	2.77	0.49
1:B:406:LYS:NZ	1:B:887:ASP:O	2.46	0.49
1:B:716:MET:HE2	1:B:878:MET:CE	2.43	0.49
1:A:645:ASN:O	1:A:648:ASN:O	2.30	0.49
1:B:329:LYS:CD	1:B:361:ASP:OD1	2.58	0.49
1:B:544:ASN:OD1	1:B:544:ASN:C	2.56	0.49
1:B:302:TYR:CE1	1:B:331:ALA:CB	2.94	0.49
1:B:667:PRO:O	1:B:670:THR:HG22	2.10	0.49
1:B:719:LEU:N	1:B:719:LEU:CD2	2.74	0.49
1:B:728:LYS:HD3	1:B:786:THR:HB	1.94	0.49
1:A:294:ARG:NH1	1:A:298:HIS:CD2	2.81	0.49
1:B:776:PRO:HB2	1:B:777:TYR:CD1	2.36	0.49
1:B:930:THR:HA	1:B:936:LEU:HD12	1.94	0.49
1:A:573:HIS:C	1:A:573:HIS:HD2	2.17	0.49
1:B:708:LEU:HD22	1:B:712:MET:HG2	1.94	0.49
1:B:716:MET:CE	1:B:878:MET:HE2	2.43	0.49
1:B:752:VAL:HG23	1:B:811:LEU:O	2.12	0.49
1:B:763:HIS:HE2	1:B:848:ALA:C	2.19	0.49
1:A:763:HIS:HD2	1:A:779:ILE:CD1	2.26	0.48
1:B:705:GLN:CG	1:B:827:ILE:HD13	2.42	0.48
1:B:734:VAL:HG13	1:B:744:LEU:HD23	1.95	0.48
1:B:930:THR:HA	1:B:936:LEU:CD1	2.43	0.48
1:B:751:THR:HG22	1:B:754:GLU:HG3	1.95	0.48
1:B:331:ALA:HB1	1:B:357:TRP:CZ2	2.43	0.48
1:B:354:LEU:C	1:B:354:LEU:CD2	2.86	0.48
1:B:362:VAL:O	1:B:365:ALA:HB3	2.14	0.48
1:B:769:HIS:ND1	1:B:815:THR:HB	2.29	0.48
1:B:826:TYR:HB3	1:B:830:ARG:O	2.13	0.48
1:B:768:LYS:HB3	1:B:769:HIS:CD2	2.49	0.48
1:A:551:PHE:CZ	1:A:555:LEU:HD11	2.49	0.48
1:B:379:ARG:HH12	1:B:407:TYR:HB2	1.77	0.48
1:B:339:ILE:HG13	1:B:340:ASN:H	1.79	0.48
1:B:841:LEU:O	1:B:932:VAL:HG21	2.14	0.48
1:A:751:THR:CG2	1:A:754:GLU:CD	2.86	0.48
1:B:812:LEU:H	1:B:812:LEU:CD2	2.26	0.48
1:A:836:PRO:HB2	1:A:837:PRO:HA	1.96	0.48
1:B:300:ILE:CG1	1:B:304:TYR:HE2	2.27	0.48
1:B:339:ILE:HG13	1:B:340:ASN:N	2.29	0.48
1:B:610:GLU:CD	1:B:643:LYS:HG3	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:631:LYS:HD2	1:B:635:LEU:HD21	1.94	0.48
1:B:807:HIS:O	1:B:810:ASN:HB2	2.14	0.48
1:A:852:ILE:HA	1:A:852:ILE:HD12	1.60	0.47
1:B:699:HIS:O	1:B:740:LYS:O	2.32	0.47
1:A:313:GLU:CD	1:A:313:GLU:N	2.70	0.47
1:A:323:PHE:HE1	1:A:349:GLN:OE1	1.96	0.47
1:A:304:TYR:CB	1:A:305:PRO:HD3	2.44	0.47
1:A:306:PRO:HD2	1:A:308:TYR:HD2	1.73	0.47
1:B:636:LEU:HD12	1:B:670:THR:HG21	1.96	0.47
1:A:561:GLU:CD	1:A:561:GLU:N	2.71	0.47
1:A:723:GLU:OE1	1:A:723:GLU:CA	2.63	0.47
1:B:375:HIS:CD2	1:B:376:PRO:HD2	2.50	0.47
1:B:631:LYS:O	1:B:635:LEU:HD23	2.14	0.47
1:A:712:MET:HE1	1:A:878:MET:HG2	1.96	0.47
1:A:841:LEU:O	1:A:932:VAL:HG21	2.14	0.47
1:A:857:HIS:NE2	1:A:861:ARG:HD2	2.30	0.47
1:B:316:ASP:O	1:B:319:TRP:HB3	2.15	0.47
1:B:763:HIS:HE1	1:B:777:TYR:CE2	2.33	0.47
1:B:806:ARG:CZ	1:B:810:ASN:ND2	2.78	0.47
1:A:320:LYS:HD2	1:A:321:PHE:HE2	1.79	0.47
1:A:546:THR:O	1:A:549:ASN:HB3	2.15	0.47
1:A:651:PRO:HA	1:A:662:ILE:O	2.14	0.47
1:A:835:MET:HG3	1:B:943:PHE:HE2	1.80	0.47
1:A:925:LEU:O	1:A:929:ILE:HG13	2.15	0.47
1:A:305:PRO:HA	1:A:308:TYR:CE2	2.50	0.47
1:B:295:ASP:O	1:B:299:THR:HG23	2.15	0.47
1:B:614:LEU:HD12	1:B:614:LEU:C	2.39	0.47
1:A:363:GLU:HB3	1:A:877:VAL:HG13	1.95	0.46
1:A:830:ARG:CZ	1:A:830:ARG:HB3	2.44	0.46
1:B:325:LEU:C	1:B:327:SER:N	2.70	0.46
1:B:575:MET:HE3	1:B:579:VAL:HG21	1.95	0.46
1:B:730:THR:O	1:B:730:THR:OG1	2.30	0.46
1:B:751:THR:HG22	1:B:754:GLU:CG	2.45	0.46
1:B:830:ARG:NH2	1:B:903:LYS:NZ	2.58	0.46
1:B:885:MET:O	1:B:888:ALA:CB	2.62	0.46
1:B:719:LEU:O	1:B:723:GLU:HG2	2.16	0.46
1:B:843:LYS:CA	1:B:932:VAL:CG1	2.92	0.46
1:A:335:PHE:CD1	1:A:357:TRP:CH2	2.87	0.46
1:A:835:MET:HG3	1:B:943:PHE:CE2	2.50	0.46
1:B:408:GLU:CB	1:B:409:ASP:HA	2.45	0.46
1:B:654:PHE:HD1	1:B:655:PRO:CD	2.27	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:598:PHE:C	1:B:598:PHE:CD2	2.94	0.46
1:B:798:THR:CG2	1:B:803:VAL:HB	2.41	0.46
1:B:407:TYR:N	1:B:407:TYR:CD2	2.78	0.46
1:B:634:LYS:HE2	1:B:634:LYS:CA	2.43	0.46
1:A:339:ILE:C	1:A:341:TRP:N	2.73	0.46
1:A:308:TYR:HB3	1:A:310:LEU:HB2	1.97	0.46
1:A:862:LYS:O	1:A:866:THR:CG2	2.64	0.46
1:B:383:VAL:HA	1:B:386:LEU:HD12	1.97	0.46
1:B:389:ALA:HA	1:B:390:PRO:HD2	1.73	0.46
1:B:575:MET:HE1	1:B:576:TYR:HD2	1.81	0.46
1:A:312:SER:O	1:A:315:GLN:HG3	2.16	0.46
1:A:610:GLU:OE1	1:A:643:LYS:HB2	2.16	0.46
1:B:326:SER:HA	1:B:357:TRP:CE2	2.51	0.46
1:B:888:ALA:O	1:B:889:THR:CG2	2.59	0.45
1:A:922:GLN:O	1:A:926:ASP:OD2	2.34	0.45
1:B:547:LEU:HD23	1:B:547:LEU:HA	1.83	0.45
1:B:612:VAL:HG22	1:B:741:HIS:HD2	1.80	0.45
1:A:306:PRO:HD3	1:A:308:TYR:HE2	1.81	0.45
1:B:777:TYR:HB2	1:B:779:ILE:CD1	2.46	0.45
1:B:763:HIS:NE2	1:B:848:ALA:O	2.41	0.45
1:A:869:LEU:HD23	1:A:869:LEU:HA	1.85	0.45
1:A:339:ILE:CG2	1:A:350:ALA:HB2	2.43	0.45
1:A:400:GLN:HG2	1:A:881:LEU:CD2	2.46	0.45
1:B:552:TYR:OH	1:B:605:ARG:NE	2.50	0.45
1:A:732:TYR:OH	1:A:822:ILE:HG12	2.16	0.45
1:A:899:LYS:HB3	1:A:903:LYS:HE2	1.98	0.45
1:A:719:LEU:O	1:A:723:GLU:HG2	2.17	0.45
1:A:782:GLU:CD	1:A:782:GLU:H	2.23	0.45
1:B:587:LEU:HB3	1:B:598:PHE:HB2	1.99	0.45
1:A:314:GLU:CA	1:A:317:LEU:HG	2.46	0.45
1:A:321:PHE:O	1:A:325:LEU:HD23	2.17	0.45
1:A:948:ARG:C	1:A:948:ARG:CD	2.88	0.45
1:A:703:LEU:HD12	1:A:742:GLY:HA3	1.99	0.44
1:B:558:GLU:HB3	1:B:576:TYR:HD1	1.79	0.44
1:B:651:PRO:HA	1:B:662:ILE:O	2.17	0.44
1:A:302:TYR:CD1	1:A:331:ALA:HA	2.50	0.44
1:A:305:PRO:HB3	1:A:308:TYR:HB2	2.00	0.44
1:A:763:HIS:HB3	1:A:767:ARG:NH2	2.32	0.44
1:B:375:HIS:CD2	1:B:376:PRO:CD	3.00	0.44
1:B:712:MET:HE1	1:B:878:MET:HG2	1.99	0.44
1:B:785:ASP:O	1:B:789:LYS:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:315:GLN:O	1:A:319:TRP:HB2	2.17	0.44
1:A:552:TYR:C	1:A:552:TYR:HD2	2.21	0.44
1:B:330:LYS:HE3	1:B:334:LYS:HE2	1.99	0.44
1:B:360:MET:SD	1:B:364:ASP:HB3	2.58	0.44
1:A:421:PHE:N	1:A:421:PHE:HD1	2.14	0.44
1:B:678:MET:HA	1:B:679:PRO:HD3	1.82	0.44
1:B:810:ASN:HA	1:B:822:ILE:CG2	2.46	0.44
1:B:830:ARG:NH1	1:B:903:LYS:NZ	2.64	0.44
1:B:403:GLN:HE22	1:B:406:LYS:HZ1	1.65	0.44
1:B:805:ASP:O	1:B:810:ASN:ND2	2.51	0.44
1:A:328:HIS:O	1:A:331:ALA:HB3	2.18	0.44
1:A:871:LEU:HD22	1:A:878:MET:HE1	2.00	0.44
1:B:575:MET:HE1	1:B:576:TYR:CD2	2.53	0.44
1:B:612:VAL:CG2	1:B:741:HIS:CD2	3.01	0.44
1:B:677:LEU:HD11	1:B:700:GLY:HA3	1.99	0.44
1:B:736:ALA:HA	1:B:742:GLY:HA3	2.00	0.44
1:A:332:LEU:O	1:A:336:LEU:HD23	2.18	0.43
1:B:342:LYS:HD2	1:B:342:LYS:HA	1.88	0.43
1:B:360:MET:CG	1:B:364:ASP:CB	2.93	0.43
1:B:379:ARG:HH12	1:B:407:TYR:CB	2.31	0.43
1:A:313:GLU:OE1	1:A:313:GLU:N	2.51	0.43
1:A:591:ASN:O	1:A:595:ARG:HG3	2.17	0.43
1:A:737:THR:HB	1:A:741:HIS:CD2	2.53	0.43
1:A:837:PRO:O	1:A:840:LYS:HE3	2.18	0.43
1:B:578:MET:O	1:B:582:MET:HG3	2.18	0.43
1:A:319:TRP:HA	1:A:322:ARG:CG	2.47	0.43
1:B:298:HIS:NE2	1:B:321:PHE:HB2	2.33	0.43
1:B:612:VAL:CG2	1:B:741:HIS:HD2	2.31	0.43
1:B:722:ARG:CA	1:B:722:ARG:NH1	2.79	0.43
1:A:292:SER:C	1:A:295:ASP:OD2	2.61	0.43
1:A:734:VAL:CG2	1:A:744:LEU:HD23	2.36	0.43
1:A:807:HIS:O	1:A:810:ASN:HB2	2.18	0.43
1:B:408:GLU:HA	1:B:410:PRO:HD3	1.99	0.43
1:B:631:LYS:CD	1:B:635:LEU:HD23	2.48	0.43
1:B:322:ARG:NH2	1:B:334:LYS:CB	2.80	0.43
1:B:631:LYS:CD	1:B:635:LEU:CD2	2.95	0.43
1:B:317:LEU:HD12	1:B:317:LEU:C	2.43	0.43
1:B:558:GLU:CB	1:B:576:TYR:CD1	3.01	0.43
1:B:660:ILE:HD12	1:B:693:TYR:CD2	2.53	0.43
1:A:553:TRP:CZ3	1:A:737:THR:HG23	2.54	0.43
1:A:723:GLU:O	1:A:724:ASN:CB	2.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:763:HIS:O	1:A:767:ARG:HB2	2.19	0.43
1:B:331:ALA:HB3	1:B:357:TRP:HE1	1.84	0.43
1:A:827:ILE:HG22	1:A:892:ASP:HB3	2.00	0.43
1:B:363:GLU:O	1:B:364:ASP:C	2.61	0.43
1:B:759:GLU:O	1:B:760:GLY:C	2.61	0.43
1:A:798:THR:CG2	1:A:803:VAL:HB	2.44	0.42
1:B:314:GLU:HB3	1:B:317:LEU:HD23	2.01	0.42
1:A:830:ARG:NH1	1:A:892:ASP:OD1	2.53	0.42
1:B:317:LEU:O	1:B:320:LYS:HG2	2.19	0.42
1:A:862:LYS:O	1:A:866:THR:HG23	2.19	0.42
1:B:398:LEU:HD13	1:B:398:LEU:C	2.44	0.42
1:A:703:LEU:HB2	1:A:736:ALA:HB2	2.01	0.42
1:A:933:MET:HA	1:A:934:PRO:HD3	1.75	0.42
1:B:618:VAL:HG21	1:B:632:PHE:CD2	2.55	0.42
1:A:362:VAL:HG13	1:A:386:LEU:HD23	2.02	0.42
1:A:552:TYR:CG	1:A:601:LEU:HD22	2.55	0.42
1:A:673:PHE:HB2	1:A:679:PRO:HD2	2.01	0.42
1:B:413:ILE:HD13	1:B:575:MET:HG3	2.00	0.42
1:A:294:ARG:CZ	1:A:298:HIS:NE2	2.83	0.42
1:A:664:LYS:HB2	1:A:685:VAL:HG22	2.00	0.42
1:A:315:GLN:HB3	1:A:338:CYS:HB3	2.01	0.42
1:A:591:ASN:ND2	1:A:591:ASN:N	2.68	0.42
1:A:762:ILE:HD13	1:A:813:LEU:HD21	2.02	0.42
1:B:575:MET:CE	1:B:576:TYR:HD2	2.33	0.42
1:B:930:THR:HG22	1:B:936:LEU:HB3	2.02	0.42
1:A:614:LEU:C	1:A:614:LEU:HD12	2.44	0.42
1:A:843:LYS:O	1:A:847:GLU:HG3	2.19	0.42
1:B:300:ILE:HA	1:B:304:TYR:CD2	2.49	0.42
1:B:719:LEU:HD12	1:B:874:HIS:CG	2.55	0.42
1:B:843:LYS:HA	1:B:932:VAL:HG11	2.01	0.42
1:A:299:THR:O	1:A:303:ARG:HB2	2.19	0.42
1:A:549:ASN:CG	1:A:658:PRO:HG3	2.45	0.42
1:B:375:HIS:CD2	1:B:376:PRO:N	2.88	0.42
1:B:763:HIS:O	1:B:767:ARG:HB2	2.20	0.42
1:B:863:GLN:C	1:B:866:THR:HG22	2.44	0.42
1:A:298:HIS:HB3	1:A:302:TYR:CE2	2.55	0.41
1:A:580:LEU:HD23	1:A:580:LEU:HA	1.86	0.41
1:A:871:LEU:HD23	1:A:871:LEU:HA	1.92	0.41
1:B:395:LEU:HD21	1:B:550:TYR:CD2	2.55	0.41
1:B:746:TYR:CE2	1:B:748:ASP:HA	2.55	0.41
1:A:648:ASN:OD1	1:A:664:LYS:HE2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:843:LYS:HA	1:A:932:VAL:HG13	2.03	0.41
1:B:305:PRO:HA	1:B:306:PRO:HD3	1.80	0.41
1:B:746:TYR:HE2	1:B:748:ASP:HA	1.85	0.41
1:A:570:GLU:O	1:A:574:ASP:OD1	2.39	0.41
1:A:677:LEU:HD22	1:A:700:GLY:CA	2.48	0.41
1:A:733:LYS:H	1:A:745:GLN:HB2	1.85	0.41
1:A:835:MET:HG3	1:A:836:PRO:HD2	2.02	0.41
1:A:872:ARG:NH1	1:A:909:GLN:C	2.79	0.41
1:A:320:LYS:HD2	1:A:321:PHE:CE2	2.56	0.41
1:A:400:GLN:HE21	1:A:885:MET:CG	2.33	0.41
1:A:638:GLU:HG2	1:A:641:MET:HB2	2.03	0.41
1:B:552:TYR:CD2	1:B:601:LEU:HD22	2.56	0.41
1:B:634:LYS:HA	1:B:634:LYS:CE	2.37	0.41
1:A:544:ASN:C	1:A:544:ASN:OD1	2.63	0.41
1:B:375:HIS:HD2	1:B:376:PRO:N	2.18	0.41
1:B:915:GLU:O	1:B:919:GLN:HG3	2.21	0.41
1:A:375:HIS:CD2	1:A:377:GLN:H	2.39	0.41
1:A:662:ILE:H	1:A:662:ILE:HG12	1.79	0.41
1:A:709:ILE:HD13	1:A:709:ILE:HA	1.90	0.41
1:A:888:ALA:C	1:A:889:THR:CG2	2.94	0.41
1:B:621:GLU:HA	1:B:622:PRO:HD3	1.90	0.41
1:B:629:THR:CG2	1:B:672:LEU:HD12	2.51	0.41
1:A:372:THR:HG1	1:A:373:PHE:HE2	1.63	0.41
1:A:405:LEU:CD2	1:A:533:LEU:HD21	2.51	0.41
1:A:533:LEU:HD23	1:A:534:CYS:N	2.36	0.41
1:B:300:ILE:CA	1:B:304:TYR:HE2	2.15	0.41
1:B:332:LEU:HD13	1:B:357:TRP:CG	2.55	0.41
1:A:634:LYS:HA	1:A:634:LYS:HD2	1.84	0.41
1:B:575:MET:HE2	1:B:576:TYR:N	2.36	0.41
1:B:591:ASN:OD1	1:B:591:ASN:C	2.64	0.41
1:A:913:THR:CG2	1:A:916:GLU:CD	2.94	0.40
1:B:783:VAL:HA	1:B:786:THR:HG23	2.03	0.40
1:B:804:GLY:O	1:B:805:ASP:HB3	2.21	0.40
1:B:556:SER:HB2	1:B:605:ARG:HH21	1.85	0.40
1:B:686:THR:OG1	1:B:691:HIS:HB2	2.21	0.40
1:B:631:LYS:HD3	1:B:635:LEU:HD23	2.03	0.40
1:B:775:GLY:O	1:B:776:PRO:C	2.64	0.40
1:A:350:ALA:O	1:A:353:MET:HB2	2.22	0.40
1:A:699:HIS:O	1:A:699:HIS:CD2	2.74	0.40
1:B:806:ARG:HA	1:B:810:ASN:HD22	1.87	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:TYR:OH	1:A:782:GLU:OE1[6_555]	2.12	0.08

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	537/696 (77%)	508 (95%)	23 (4%)	6 (1%)	11	36
1	B	537/696 (77%)	512 (95%)	21 (4%)	4 (1%)	18	48
All	All	1074/1392 (77%)	1020 (95%)	44 (4%)	10 (1%)	14	41

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	PRO
1	A	737	THR
1	B	836	PRO
1	A	304	TYR
1	B	305	PRO
1	A	836	PRO
1	B	306	PRO
1	B	408	GLU
1	A	362	VAL
1	A	776	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	489/612 (80%)	392 (80%)	97 (20%)	1	5
1	B	488/612 (80%)	408 (84%)	80 (16%)	2	8
All	All	977/1224 (80%)	800 (82%)	177 (18%)	2	6

All (177) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	293	ILE
1	A	297	LEU
1	A	303	ARG
1	A	309	VAL
1	A	313	GLU
1	A	317	LEU
1	A	333	THR
1	A	335	PHE
1	A	349	GLN
1	A	351	LEU
1	A	362	VAL
1	A	363	GLU
1	A	370	SER
1	A	373	PHE
1	A	374	THR
1	A	378	VAL
1	A	391	ASP
1	A	393	ASP
1	A	396	LEU
1	A	412	HIS
1	A	416	LEU
1	A	417	HIS
1	A	421	PHE
1	A	532	ASN
1	A	534	CYS
1	A	540	ARG
1	A	544	ASN
1	A	546	THR
1	A	549	ASN
1	A	557	ILE
1	A	561	GLU
1	A	573	HIS
1	A	580	LEU
1	A	581	LYS
1	A	591	ASN

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Mol	Chain	Res	Type
1	A	594	LEU
1	A	602	ARG
1	A	606	ARG
1	A	611	LEU
1	A	614	LEU
1	A	629	THR
1	A	634	LYS
1	A	659	GLU
1	A	662	ILE
1	A	669	ARG
1	A	671	SER
1	A	674	LYS
1	A	685	VAL
1	A	686	THR
1	A	691	HIS
1	A	708	LEU
1	A	710	LEU
1	A	711	GLN
1	A	715	LEU
1	A	723	GLU
1	A	724	ASN
1	A	725	LEU
1	A	728	LYS
1	A	730	THR
1	A	740	LYS
1	A	747	VAL
1	A	748	ASP
1	A	750	CYS
1	A	751	THR
1	A	759	GLU
1	A	786	THR
1	A	788	ILE
1	A	791	CYS
1	A	795	CYS
1	A	806	ARG
1	A	807	HIS
1	A	808	LEU
1	A	811	LEU
1	A	812	LEU
1	A	813	LEU
1	A	819	LEU
1	A	822	ILE

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Mol	Chain	Res	Type
1	A	827	ILE
1	A	831	ASP
1	A	835	MET
1	A	839	MET
1	A	842	SER
1	A	852	ILE
1	A	854	SER
1	A	855	GLU
1	A	866	THR
1	A	877	VAL
1	A	889	THR
1	A	896	GLU
1	A	902	LYS
1	A	909	GLN
1	A	913	THR
1	A	927	VAL
1	A	930	THR
1	A	932	VAL
1	A	940	ILE
1	A	942	ARG
1	B	298	HIS
1	B	299	THR
1	B	300	ILE
1	B	313	GLU
1	B	314	GLU
1	B	316	ASP
1	B	317	LEU
1	B	318	VAL
1	B	321	PHE
1	B	333	THR
1	B	338	CYS
1	B	343	LEU
1	B	348	THR
1	B	351	LEU
1	B	354	LEU
1	B	361	ASP
1	B	368	LEU
1	B	372	THR
1	B	374	THR
1	B	394	LEU
1	B	398	LEU
1	B	399	LEU

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Mol	Chain	Res	Type
1	B	414	VAL
1	B	549	ASN
1	B	555	LEU
1	B	571	ARG
1	B	573	HIS
1	B	575	MET
1	B	580	LEU
1	B	592	PHE
1	B	594	LEU
1	B	605	ARG
1	B	611	LEU
1	B	614	LEU
1	B	634	LYS
1	B	645	ASN
1	B	656	LEU
1	B	660	ILE
1	B	663	THR
1	B	670	THR
1	B	677	LEU
1	B	682	LEU
1	B	685	VAL
1	B	699	HIS
1	B	702	ASP
1	B	708	LEU
1	B	710	LEU
1	B	714	THR
1	B	715	LEU
1	B	719	LEU
1	B	722	ARG
1	B	725	LEU
1	B	730	THR
1	B	741	HIS
1	B	746	TYR
1	B	747	VAL
1	B	750	CYS
1	B	751	THR
1	B	758	ARG
1	B	759	GLU
1	B	782	GLU
1	B	786	THR
1	B	788	ILE
1	B	811	LEU

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Mol	Chain	Res	Type
1	B	812	LEU
1	B	813	LEU
1	B	819	LEU
1	B	822	ILE
1	B	827	ILE
1	B	839	MET
1	B	842	SER
1	B	852	ILE
1	B	876	ASN
1	B	889	THR
1	B	898	ASP
1	B	915	GLU
1	B	923	SER
1	B	927	VAL
1	B	937	VAL
1	B	945	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	349	GLN
1	A	568	GLN
1	A	573	HIS
1	A	591	ASN
1	A	639	GLN
1	A	645	ASN
1	A	699	HIS
1	A	711	GLN
1	A	724	ASN
1	A	741	HIS
1	A	763	HIS
1	A	870	HIS
1	A	876	ASN
1	A	922	GLN
1	A	939	GLN
1	B	340	ASN
1	B	349	GLN
1	B	375	HIS
1	B	403	GLN
1	B	633	GLN
1	B	639	GLN
1	B	691	HIS

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Mol	Chain	Res	Type
1	B	711	GLN
1	B	724	ASN
1	B	741	HIS
1	B	745	GLN
1	B	807	HIS
1	B	810	ASN
1	B	821	HIS
1	B	876	ASN
1	B	919	GLN
1	B	922	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	543/696 (78%)	0.57	35 (6%) 25 20	20, 33, 61, 86	0
1	B	543/696 (78%)	0.30	25 (4%) 37 29	22, 40, 74, 87	0
All	All	1086/1392 (78%)	0.43	60 (5%) 30 24	20, 36, 70, 87	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	302	TYR	9.6
1	B	322	ARG	8.6
1	B	302	TYR	6.5
1	B	420	ILE	6.2
1	A	294	ARG	6.1
1	A	322	ARG	6.0
1	B	306	PRO	5.9
1	A	341	TRP	5.6
1	A	324	TYR	5.4
1	B	298	HIS	4.4
1	B	293	ILE	4.3
1	A	323	PHE	4.3
1	B	324	TYR	4.2
1	A	311	SER	4.2
1	A	310	LEU	3.8
1	A	340	ASN	3.8
1	A	293	ILE	3.7
1	B	325	LEU	3.7
1	A	325	LEU	3.6
1	A	306	PRO	3.5
1	B	291	ALA	3.5
1	A	328	HIS	3.3
1	B	343	LEU	3.3
1	A	291	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	340	ASN	3.2
1	B	341	TRP	3.2
1	B	310	LEU	3.2
1	B	307	THR	3.2
1	B	305	PRO	2.9
1	A	303	ARG	2.9
1	B	297	LEU	2.8
1	B	677	LEU	2.7
1	A	298	HIS	2.6
1	B	319	TRP	2.6
1	B	342	LYS	2.6
1	B	321	PHE	2.5
1	A	948	ARG	2.5
1	B	362	VAL	2.4
1	A	352	TRP	2.4
1	A	292	SER	2.3
1	A	308	TYR	2.3
1	B	690	HIS	2.3
1	A	339	ILE	2.3
1	A	326	SER	2.3
1	A	376	PRO	2.3
1	A	759	GLU	2.3
1	B	303	ARG	2.3
1	A	401	LEU	2.2
1	A	906	GLU	2.2
1	A	567	LYS	2.2
1	A	722	ARG	2.2
1	B	416	LEU	2.2
1	A	377	GLN	2.2
1	A	410	PRO	2.2
1	B	317	LEU	2.2
1	A	373	PHE	2.1
1	A	369	LEU	2.1
1	A	305	PRO	2.0
1	A	297	LEU	2.0
1	A	408	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.