



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:10 PM UTC

PDB ID : 4AV3 / pdb_00004av3
Title : Crystal structure of Thermotoga Maritima sodium pumping membrane integral pyrophosphatase with metal ions in active site
Authors : Kajander, T.; Kogan, K.; Kellosalo, J.; Pokharel, K.; Goldman, A.
Deposited on : 2012-05-23
Resolution : 2.60 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

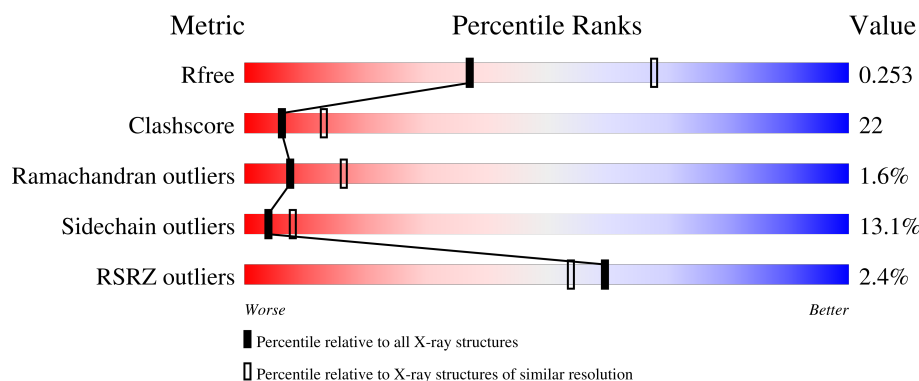
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	735	2% 64% 23% 8% 6%
1	B	735	2% 64% 25% 7% ...

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9988 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 K(+)-STIMULATED PYROPHOSPHATE-ENERGIZED SODIUM PUMP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	694	Total	C	N	O	S	0	0	0
			4926	3230	769	900	27			
1	B	707	Total	C	N	O	S	0	0	0
			5039	3294	789	929	27			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	MET	-	expression tag	UNP Q9S5X0
A	-7	ARG	-	expression tag	UNP Q9S5X0
A	-6	GLY	-	expression tag	UNP Q9S5X0
A	-5	SER	-	expression tag	UNP Q9S5X0
A	-4	HIS	-	expression tag	UNP Q9S5X0
A	-3	HIS	-	expression tag	UNP Q9S5X0
A	-2	HIS	-	expression tag	UNP Q9S5X0
A	-1	HIS	-	expression tag	UNP Q9S5X0
A	0	HIS	-	expression tag	UNP Q9S5X0
A	1	HIS	-	expression tag	UNP Q9S5X0
A	353	LEU	VAL	engineered mutation	UNP Q9S5X0
A	395	GLY	SER	engineered mutation	UNP Q9S5X0
B	-8	MET	-	expression tag	UNP Q9S5X0
B	-7	ARG	-	expression tag	UNP Q9S5X0
B	-6	GLY	-	expression tag	UNP Q9S5X0
B	-5	SER	-	expression tag	UNP Q9S5X0
B	-4	HIS	-	expression tag	UNP Q9S5X0
B	-3	HIS	-	expression tag	UNP Q9S5X0
B	-2	HIS	-	expression tag	UNP Q9S5X0
B	-1	HIS	-	expression tag	UNP Q9S5X0
B	0	HIS	-	expression tag	UNP Q9S5X0
B	1	HIS	-	expression tag	UNP Q9S5X0
B	353	LEU	VAL	engineered mutation	UNP Q9S5X0
B	395	GLY	SER	engineered mutation	UNP Q9S5X0

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

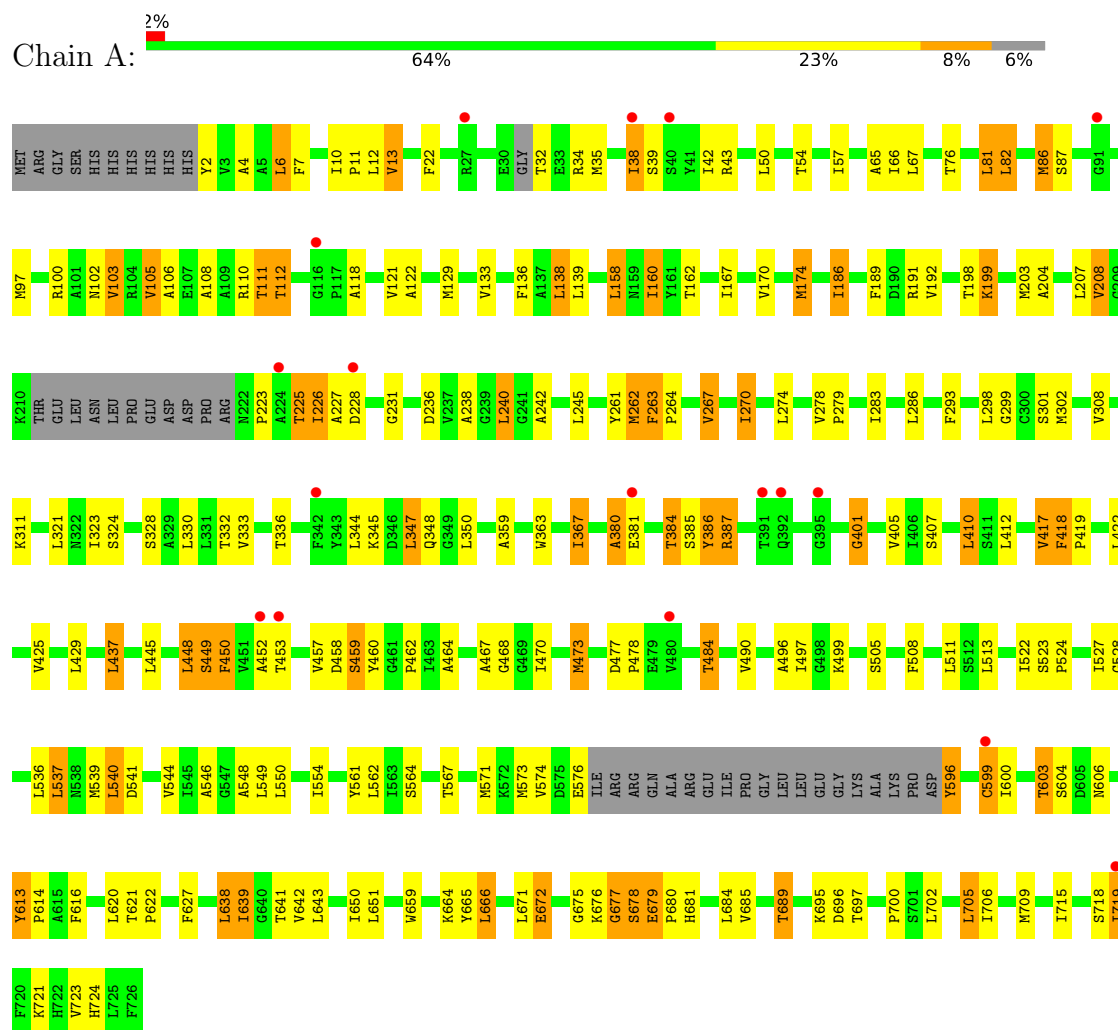
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total	O	0	0
			11	11		
4	B	8	Total	O	0	0
			8	8		

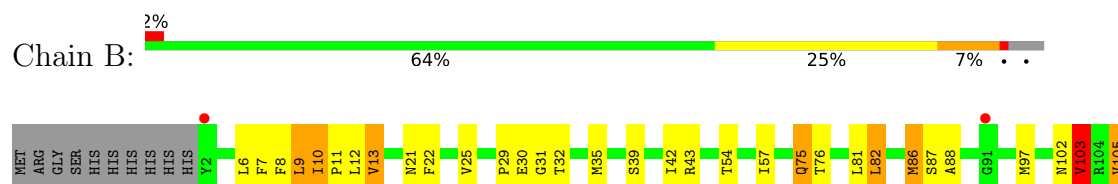
3 Residue-property plots

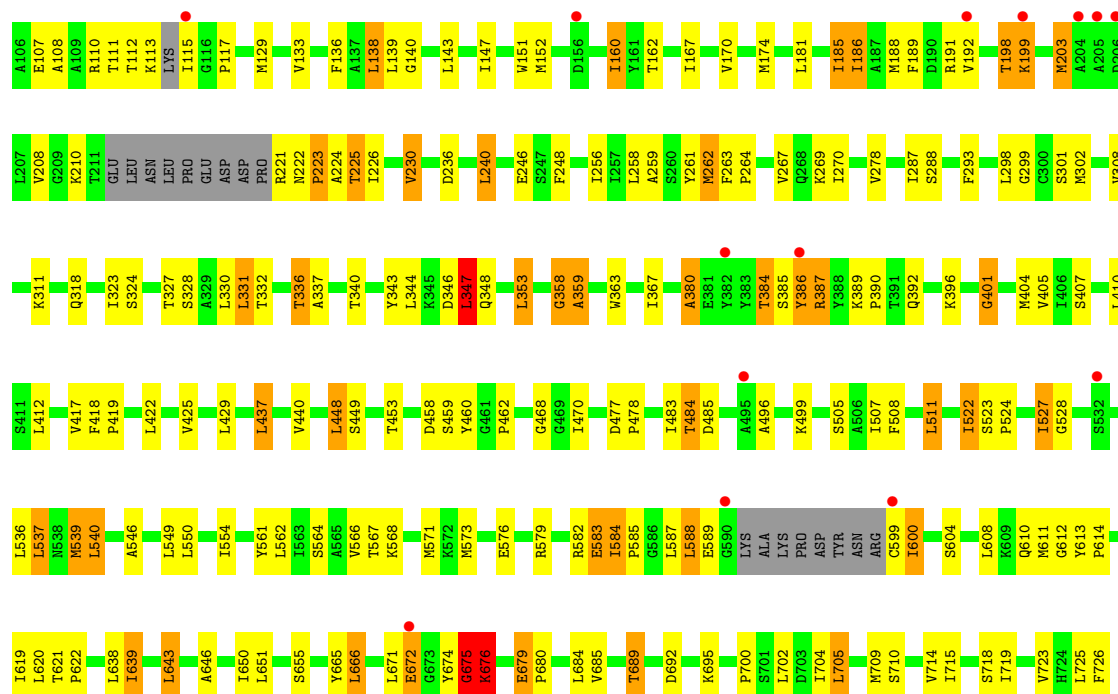
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: K(+)-STIMULATED PYROPHOSPHATE-ENERGIZED SODIUM PUMP



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.52Å 107.78Å 102.52Å 90.00° 108.50° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 50.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.8 (50.00-2.60) 97.8 (50.00-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.61Å)	Xtriage
Refinement program		Depositor
R, R_{free}	0.205 , 0.245 (Not available) , 0.253	Depositor DCC
R_{free} test set	2622 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	71.0	Xtriage
Anisotropy	0.654	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 81.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9988	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MG

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	0.99	2/5031 (0.0%)	1.15	22/6867 (0.3%)
1	B	0.97	3/5142 (0.1%)	1.11	14/7011 (0.2%)
All	All	0.98	5/10173 (0.0%)	1.13	36/13878 (0.3%)

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	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	452	ALA	N-CA	5.87	1.53	1.46
1	B	539	MET	N-CA	-5.67	1.38	1.46
1	B	258	LEU	CA-C	5.45	1.59	1.52
1	B	511	LEU	CA-C	5.38	1.59	1.52
1	A	238	ALA	CA-C	5.30	1.59	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	386	TYR	N-CA-C	-8.94	98.62	110.24
1	A	677	GLY	N-CA-C	-8.24	98.84	111.12
1	A	724	HIS	N-CA-C	8.09	121.03	110.43
1	B	346	ASP	N-CA-C	7.46	120.43	110.88
1	B	401	GLY	N-CA-C	7.26	119.54	111.85
1	B	449	SER	N-CA-C	7.20	121.16	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	380	ALA	N-CA-C	-6.82	103.93	111.36
1	B	539	MET	CG-SD-CE	-6.65	86.28	100.90
1	A	449	SER	N-CA-C	-6.44	97.07	110.80
1	A	160	ILE	N-CA-CB	-6.27	103.60	111.31
1	A	401	GLY	N-CA-C	6.09	118.30	111.85
1	A	450	PHE	N-CA-C	5.95	123.46	110.80
1	B	105	VAL	N-CA-C	-5.78	105.48	110.74
1	B	386	TYR	N-CA-C	-5.70	98.66	110.80
1	A	418	PHE	CA-C-N	-5.65	114.56	120.38
1	A	418	PHE	C-N-CA	-5.65	114.56	120.38
1	B	676	LYS	N-CA-C	5.53	122.58	110.80
1	A	174	MET	CG-SD-CE	-5.51	88.77	100.90
1	B	10	ILE	CB-CA-C	-5.46	108.56	114.35
1	A	279	PRO	N-CA-C	5.44	119.84	111.57
1	A	242	ALA	N-CA-C	-5.39	105.43	112.23
1	A	721	LYS	N-CA-C	-5.38	105.42	111.28
1	A	347	LEU	N-CA-C	5.35	118.27	111.69
1	B	347	LEU	N-CA-C	5.31	118.23	111.69
1	B	103	VAL	CB-CA-C	-5.30	105.09	111.88
1	A	66	ILE	N-CA-CB	5.30	117.74	110.54
1	A	278	VAL	CA-C-N	5.21	125.41	120.31
1	A	278	VAL	C-N-CA	5.21	125.41	120.31
1	A	330	LEU	N-CA-C	-5.20	105.50	111.07
1	B	440	VAL	N-CA-C	-5.11	105.56	110.72
1	A	345	LYS	N-CA-C	5.10	117.50	111.33
1	B	380	ALA	N-CA-C	-5.08	105.45	111.69
1	B	330	LEU	N-CA-C	-5.06	105.66	111.07
1	B	675	GLY	N-CA-C	5.05	125.15	113.18
1	A	263	PHE	CA-C-N	-5.03	113.86	119.19
1	A	263	PHE	C-N-CA	-5.03	113.86	119.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	528	GLY	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	4926	0	4792	212	0
1	B	5039	0	4932	240	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	11	0	0	0	0
4	B	8	0	0	2	0
All	All	9988	0	9724	442	0

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 22.

All (442) 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:B:108:ALA:O	1:B:112:THR:HG22	1.21	1.31
1:A:105:VAL:HG21	1:A:467:ALA:CB	1.83	1.08
1:A:105:VAL:HG21	1:A:467:ALA:HB1	1.11	1.07
1:A:105:VAL:CG2	1:A:467:ALA:HB1	1.84	1.07
1:A:336:THR:HG22	1:A:363:TRP:HD1	1.17	1.06
1:A:223:PRO:HG2	1:A:225:THR:HG22	1.33	1.05
1:B:336:THR:HG22	1:B:363:TRP:HD1	1.19	1.04
1:B:108:ALA:O	1:B:112:THR:CG2	2.07	1.02
1:B:584:ILE:HD13	1:B:588:LEU:HD13	1.37	1.01
1:A:336:THR:HG22	1:A:363:TRP:CD1	1.97	0.99
1:B:363:TRP:CZ2	1:B:367:ILE:HD11	1.99	0.98
1:B:336:THR:HG22	1:B:363:TRP:CD1	1.99	0.97
1:B:524:PRO:O	1:B:527:ILE:HD12	1.67	0.94
1:B:363:TRP:CE2	1:B:367:ILE:HD11	2.03	0.93
1:B:160:ILE:HD12	1:B:160:ILE:H	1.34	0.92
1:A:638:LEU:O	1:A:642:VAL:HG23	1.70	0.91
1:B:719:ILE:HG22	1:B:723:VAL:HG11	1.50	0.90
1:B:468:GLY:HA2	1:B:484:THR:HG23	1.53	0.90
1:B:719:ILE:O	1:B:723:VAL:HG12	1.73	0.89
1:A:719:ILE:HG22	1:A:723:VAL:HG11	1.54	0.89
1:B:223:PRO:HD2	1:B:225:THR:CG2	2.04	0.88
1:A:719:ILE:HG22	1:A:723:VAL:CG1	2.03	0.88
1:A:22:PHE:HA	1:A:97:MET:CE	2.05	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:MET:HE1	1:B:715:ILE:HG23	1.56	0.86
1:B:262:MET:HE3	1:B:353:LEU:HD21	1.55	0.86
1:A:86:MET:HE1	1:A:139:LEU:HD23	1.58	0.85
1:A:22:PHE:HA	1:A:97:MET:HE1	1.57	0.85
1:A:105:VAL:CG2	1:A:106:ALA:N	2.39	0.85
1:A:540:LEU:HD22	1:B:437:LEU:HD21	1.58	0.85
1:B:223:PRO:HD2	1:B:225:THR:HG22	1.58	0.83
1:A:118:ALA:O	1:A:121:VAL:HG22	1.78	0.83
1:A:468:GLY:HA2	1:A:484:THR:HG23	1.61	0.82
1:A:386:TYR:O	1:A:387:ARG:CB	2.28	0.81
1:A:174:MET:HE1	1:A:715:ILE:HG12	1.63	0.81
1:A:363:TRP:CZ2	1:A:367:ILE:HD11	2.17	0.80
1:B:386:TYR:O	1:B:387:ARG:CB	2.30	0.79
1:B:185:ILE:HD12	1:B:185:ILE:C	2.07	0.79
1:B:719:ILE:HG22	1:B:723:VAL:CG1	2.13	0.78
1:A:105:VAL:HG22	1:A:106:ALA:N	1.98	0.78
1:B:188:MET:HE1	1:B:619:ILE:HD11	1.66	0.78
1:A:34:ARG:O	1:A:38:ILE:HG22	1.83	0.78
1:B:86:MET:HE1	1:B:139:LEU:HD23	1.66	0.77
1:B:337:ALA:HB2	1:B:363:TRP:CE2	2.20	0.77
1:B:259:ALA:HA	1:B:262:MET:HE2	1.66	0.77
1:A:270:ILE:HG23	1:A:270:ILE:O	1.85	0.77
1:A:539:MET:HE3	1:B:539:MET:HE3	1.66	0.76
1:B:7:PHE:HA	1:B:10:ILE:HD13	1.67	0.76
1:A:223:PRO:HG2	1:A:225:THR:CG2	2.13	0.76
1:B:672:GLU:OE1	1:B:672:GLU:CA	2.35	0.75
1:B:672:GLU:OE1	1:B:672:GLU:HA	1.86	0.75
1:B:468:GLY:HA2	1:B:484:THR:CG2	2.16	0.75
1:A:311:LYS:HE3	1:A:323:ILE:HD12	1.69	0.74
1:A:158:LEU:HD23	1:A:263:PHE:CG	2.24	0.73
1:B:386:TYR:O	1:B:387:ARG:HB3	1.86	0.73
1:A:336:THR:CG2	1:A:363:TRP:HD1	1.98	0.73
1:A:225:THR:O	1:A:228:ASP:N	2.22	0.73
1:B:336:THR:CG2	1:B:363:TRP:HD1	2.01	0.72
1:B:584:ILE:HG23	1:B:588:LEU:HB2	1.72	0.72
1:B:725:LEU:HD13	1:B:725:LEU:C	2.16	0.71
1:B:75:GLN:H	1:B:75:GLN:HE21	1.37	0.70
1:B:240:LEU:HD13	1:B:458:ASP:HB3	1.71	0.70
1:B:270:ILE:HG22	4:B:2004:HOH:O	1.90	0.70
1:B:674:TYR:O	1:B:676:LYS:N	2.25	0.70
1:A:293:PHE:CE2	1:A:336:THR:HG23	2.27	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:PRO:CG	1:A:225:THR:HG22	2.18	0.69
1:A:540:LEU:HD22	1:B:437:LEU:CD2	2.21	0.69
1:B:599:CYS:SG	1:B:600:ILE:N	2.65	0.69
1:B:226:ILE:O	1:B:230:VAL:HG13	1.91	0.69
1:B:380:ALA:O	1:B:384:THR:HB	1.92	0.69
1:B:203:MET:CE	1:B:567:THR:HG22	2.23	0.69
1:B:705:LEU:CD2	1:B:709:MET:HE3	2.23	0.69
1:B:358:GLY:O	1:B:359:ALA:HB3	1.94	0.68
1:B:170:VAL:HG23	1:B:261:TYR:CE1	2.29	0.68
1:A:207:LEU:O	1:A:208:VAL:HG13	1.94	0.68
1:B:86:MET:CE	1:B:139:LEU:HD23	2.23	0.67
1:A:105:VAL:HG22	1:A:106:ALA:H	1.58	0.67
1:B:42:ILE:HD12	1:B:43:ARG:N	2.09	0.67
1:A:567:THR:HB	1:B:404:MET:HE1	1.77	0.67
1:B:725:LEU:HD12	1:B:726:PHE:CG	2.30	0.66
1:A:536:LEU:HD12	1:A:537:LEU:N	2.11	0.66
1:B:328:SER:CB	1:B:453:THR:HG21	2.25	0.66
1:A:311:LYS:HE3	1:A:323:ILE:CD1	2.26	0.66
1:B:138:LEU:HD13	1:B:248:PHE:CE2	2.29	0.66
1:B:340:THR:HG23	1:B:344:LEU:HD12	1.76	0.66
1:B:208:VAL:HG22	1:B:210:LYS:H	1.60	0.66
1:B:86:MET:HE3	1:B:139:LEU:HB3	1.77	0.65
1:B:138:LEU:HD13	1:B:248:PHE:HE2	1.60	0.65
1:A:723:VAL:HG22	1:A:723:VAL:O	1.96	0.65
1:A:321:LEU:HD13	1:A:457:VAL:HG13	1.79	0.65
1:B:293:PHE:CE2	1:B:336:THR:HG23	2.31	0.65
1:A:705:LEU:CD2	1:A:709:MET:HE3	2.26	0.65
1:A:600:ILE:O	1:A:604:SER:HB2	1.96	0.64
1:A:293:PHE:CE2	1:A:336:THR:CG2	2.80	0.64
1:B:363:TRP:CH2	1:B:367:ILE:HD11	2.32	0.64
1:A:328:SER:CB	1:A:453:THR:HG21	2.27	0.64
1:A:32:THR:HG22	1:A:34:ARG:H	1.63	0.63
1:B:353:LEU:O	1:B:353:LEU:HD13	1.98	0.63
1:B:293:PHE:CE2	1:B:336:THR:CG2	2.81	0.63
1:B:527:ILE:HD13	1:B:528:GLY:H	1.62	0.63
1:A:380:ALA:O	1:A:384:THR:HB	1.98	0.63
1:B:278:VAL:HG11	1:B:353:LEU:HD23	1.81	0.63
1:A:363:TRP:CZ2	1:A:367:ILE:CD1	2.81	0.63
1:B:725:LEU:HD12	1:B:726:PHE:CD2	2.33	0.63
1:A:677:GLY:O	1:A:678:SER:OG	2.16	0.63
1:A:384:THR:HG21	1:A:496:ALA:HB1	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:613:TYR:HB3	1:A:614:PRO:HD3	1.81	0.62
1:B:600:ILE:O	1:B:604:SER:HB2	1.98	0.62
1:A:468:GLY:HA2	1:A:484:THR:CG2	2.28	0.62
1:B:39:SER:HB2	1:B:103:VAL:HG13	1.80	0.62
1:B:468:GLY:CA	1:B:484:THR:HG23	2.26	0.62
1:A:86:MET:HE3	1:A:136:PHE:HD1	1.65	0.62
1:A:705:LEU:HD21	1:A:709:MET:HE3	1.80	0.62
1:A:719:ILE:HG22	1:A:723:VAL:HG12	1.82	0.61
1:B:344:LEU:HA	1:B:347:LEU:HD21	1.81	0.61
1:A:270:ILE:O	1:A:270:ILE:CG2	2.48	0.61
1:B:192:VAL:HG11	1:B:614:PRO:HB2	1.81	0.61
1:A:651:LEU:HD23	1:A:651:LEU:C	2.26	0.61
1:B:31:GLY:HA3	1:B:35:MET:HE3	1.81	0.60
1:A:264:PRO:O	1:A:267:VAL:HG13	2.01	0.60
1:B:262:MET:CE	1:B:353:LEU:HD21	2.29	0.60
1:B:160:ILE:H	1:B:160:ILE:CD1	2.06	0.60
1:B:536:LEU:HD12	1:B:537:LEU:N	2.16	0.60
1:B:651:LEU:C	1:B:651:LEU:HD23	2.27	0.60
1:A:105:VAL:HA	1:A:121:VAL:HG21	1.82	0.60
1:A:385:SER:C	1:A:386:TYR:O	2.43	0.60
1:B:22:PHE:HA	1:B:97:MET:CE	2.32	0.59
1:A:121:VAL:HG23	1:A:122:ALA:N	2.16	0.59
1:A:616:PHE:CD2	1:A:620:LEU:CD1	2.85	0.59
1:B:705:LEU:HD21	1:B:709:MET:HE3	1.85	0.59
1:B:167:ILE:HD13	1:B:719:ILE:HG12	1.85	0.59
1:A:2:TYR:CZ	1:A:4:ALA:HB2	2.38	0.58
1:A:86:MET:HE2	1:A:139:LEU:HB3	1.84	0.58
1:A:39:SER:HB2	1:A:103:VAL:HG13	1.84	0.58
1:A:105:VAL:O	1:A:108:ALA:HB3	2.02	0.58
1:B:666:LEU:HB2	1:B:671:LEU:HD12	1.85	0.58
1:A:199:LYS:HD3	1:A:696:ASP:HB2	1.85	0.58
1:A:308:VAL:HG21	1:A:324:SER:HB3	1.85	0.58
1:A:464:ALA:O	1:A:467:ALA:HB3	2.03	0.58
1:B:685:VAL:O	1:B:689:THR:HG23	2.04	0.58
1:B:705:LEU:HD21	1:B:709:MET:CE	2.33	0.57
1:A:38:ILE:HD13	1:A:470:ILE:HA	1.85	0.57
1:B:112:THR:O	1:B:113:LYS:CB	2.52	0.57
1:A:204:ALA:HB3	1:A:573:MET:HE2	1.86	0.57
1:A:86:MET:CE	1:A:139:LEU:HD23	2.31	0.57
1:A:12:LEU:CD1	1:A:139:LEU:HD13	2.35	0.57
1:A:381:GLU:OE1	1:A:497:ILE:HD11	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:ALA:HB1	1:B:425:VAL:HG13	1.86	0.56
1:A:616:PHE:CD2	1:A:620:LEU:HD12	2.39	0.56
1:A:199:LYS:HD3	1:A:696:ASP:CB	2.35	0.56
1:A:299:GLY:HA2	1:A:302:MET:HE2	1.87	0.56
1:A:596:TYR:N	1:A:596:TYR:CD1	2.73	0.56
1:A:599:CYS:O	1:A:603:THR:HG22	2.06	0.56
1:B:174:MET:HE1	1:B:715:ILE:CG2	2.33	0.56
1:B:384:THR:HG21	1:B:496:ALA:HB1	1.87	0.56
1:A:192:VAL:HG13	1:A:562:LEU:HD21	1.87	0.56
1:B:562:LEU:O	1:B:566:VAL:HG23	2.06	0.56
1:A:223:PRO:C	1:A:225:THR:H	2.12	0.56
1:A:204:ALA:CB	1:A:573:MET:HE2	2.35	0.56
1:B:608:LEU:HD23	1:B:611:MET:HE2	1.88	0.56
1:A:42:ILE:HD12	1:A:43:ARG:N	2.21	0.56
1:B:22:PHE:HA	1:B:97:MET:HE1	1.86	0.55
1:B:385:SER:C	1:B:386:TYR:O	2.40	0.55
1:A:86:MET:HE3	1:A:136:PHE:HA	1.89	0.55
1:B:448:LEU:HD21	1:B:505:SER:HB3	1.88	0.55
1:A:105:VAL:CG2	1:A:467:ALA:CB	2.61	0.55
1:A:459:SER:O	1:A:462:PRO:HD2	2.07	0.55
1:A:448:LEU:C	1:A:449:SER:O	2.47	0.55
1:A:468:GLY:CA	1:A:484:THR:HG23	2.35	0.55
1:A:573:MET:O	1:A:576:GLU:N	2.39	0.55
1:B:705:LEU:HD22	1:B:709:MET:HE3	1.88	0.55
1:A:105:VAL:HG23	1:A:106:ALA:N	2.18	0.55
1:B:223:PRO:HG2	1:B:224:ALA:H	1.72	0.54
1:B:358:GLY:O	1:B:359:ALA:CB	2.53	0.54
1:A:719:ILE:CG2	1:A:723:VAL:HG11	2.31	0.54
1:B:459:SER:O	1:B:462:PRO:HD2	2.07	0.54
1:B:725:LEU:C	1:B:725:LEU:CD1	2.81	0.54
1:A:719:ILE:O	1:A:723:VAL:HG12	2.07	0.54
1:A:32:THR:HB	1:A:35:MET:HG3	1.88	0.54
1:B:418:PHE:HB3	1:B:419:PRO:HD3	1.89	0.54
1:B:21:ASN:O	1:B:25:VAL:HG23	2.08	0.54
1:B:675:GLY:O	1:B:676:LYS:CB	2.55	0.54
1:B:412:LEU:C	1:B:412:LEU:HD23	2.33	0.54
1:B:167:ILE:CG2	1:B:718:SER:HB3	2.38	0.54
1:A:344:LEU:HD22	1:A:347:LEU:CD1	2.37	0.53
1:A:437:LEU:HD21	1:B:540:LEU:HD22	1.90	0.53
1:A:118:ALA:O	1:A:121:VAL:CG2	2.52	0.53
1:A:344:LEU:CD2	1:A:347:LEU:CD1	2.87	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:549:LEU:HD11	1:B:511:LEU:HD11	1.90	0.53
1:A:225:THR:O	1:A:226:ILE:C	2.51	0.53
1:B:666:LEU:C	1:B:666:LEU:HD12	2.33	0.53
1:A:167:ILE:CG2	1:A:718:SER:HB3	2.39	0.53
1:B:584:ILE:O	1:B:587:LEU:N	2.40	0.53
1:B:672:GLU:OE1	1:B:672:GLU:N	2.41	0.53
1:A:336:THR:CG2	1:A:363:TRP:CD1	2.82	0.53
1:A:671:LEU:O	1:A:672:GLU:CB	2.56	0.53
1:B:108:ALA:C	1:B:112:THR:HG22	2.20	0.53
1:A:418:PHE:HB3	1:A:419:PRO:HD3	1.91	0.53
1:B:527:ILE:HD13	1:B:528:GLY:N	2.24	0.53
1:B:613:TYR:HB3	1:B:614:PRO:HD3	1.91	0.53
1:B:160:ILE:HD12	1:B:160:ILE:N	2.13	0.53
1:B:185:ILE:HD12	1:B:185:ILE:O	2.09	0.53
1:B:621:THR:HB	1:B:622:PRO:HD3	1.90	0.53
1:A:203:MET:SD	1:A:567:THR:HG22	2.49	0.53
1:B:308:VAL:HG21	1:B:324:SER:HB3	1.90	0.53
1:A:548:ALA:HB1	1:B:643:LEU:HD21	1.91	0.52
1:B:112:THR:HG21	1:B:117:PRO:HG2	1.89	0.52
1:B:522:ILE:HD13	1:B:523:SER:N	2.24	0.52
1:A:448:LEU:HD21	1:A:505:SER:HB3	1.92	0.52
1:A:536:LEU:HD12	1:A:536:LEU:C	2.35	0.52
1:B:224:ALA:HB1	1:B:573:MET:HE2	1.91	0.52
1:B:224:ALA:HB1	1:B:573:MET:CE	2.39	0.52
1:A:39:SER:OG	1:A:103:VAL:HG22	2.10	0.52
1:A:677:GLY:O	1:A:678:SER:CB	2.58	0.52
1:B:723:VAL:O	1:B:723:VAL:HG22	2.10	0.52
1:A:384:THR:HG21	1:A:496:ALA:CB	2.40	0.52
1:B:76:THR:HG21	1:B:174:MET:HE2	1.92	0.51
1:B:138:LEU:CD1	1:B:248:PHE:CE2	2.92	0.51
1:A:42:ILE:HD12	1:A:42:ILE:C	2.35	0.51
1:B:203:MET:HE1	1:B:567:THR:HG22	1.91	0.51
1:B:299:GLY:HA2	1:B:302:MET:HE2	1.93	0.51
1:A:344:LEU:HD23	1:A:347:LEU:HD11	1.91	0.51
1:A:676:LYS:HA	1:A:681:HIS:CE1	2.45	0.51
1:B:129:MET:HE3	1:B:133:VAL:CG2	2.40	0.51
1:A:158:LEU:HD23	1:A:263:PHE:CD2	2.46	0.51
1:B:9:LEU:HD23	1:B:9:LEU:N	2.26	0.51
1:A:561:TYR:C	1:A:561:TYR:CD1	2.88	0.51
1:A:621:THR:HB	1:A:622:PRO:HD3	1.93	0.51
1:A:675:GLY:C	1:A:677:GLY:H	2.19	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:725:LEU:HD12	1:B:726:PHE:CD1	2.45	0.50
1:B:39:SER:HB2	1:B:103:VAL:CG1	2.41	0.50
1:A:262:MET:HE1	1:A:286:LEU:HD13	1.91	0.50
1:A:666:LEU:C	1:A:666:LEU:HD12	2.36	0.50
1:B:129:MET:HE3	1:B:133:VAL:HG23	1.93	0.50
1:B:344:LEU:HA	1:B:347:LEU:CD2	2.41	0.50
1:B:7:PHE:HA	1:B:10:ILE:CD1	2.39	0.50
1:B:499:LYS:CE	1:B:695:LYS:O	2.59	0.50
1:A:22:PHE:HA	1:A:97:MET:HE2	1.90	0.50
1:A:42:ILE:HD11	1:A:102:ASN:HD21	1.76	0.50
1:B:181:LEU:O	1:B:185:ILE:HG23	2.12	0.50
1:B:460:TYR:CD1	1:B:460:TYR:C	2.89	0.50
1:B:29:PRO:C	1:B:107:GLU:OE1	2.54	0.50
1:B:31:GLY:CA	1:B:35:MET:HE3	2.40	0.50
1:B:311:LYS:HE3	1:B:323:ILE:HD12	1.94	0.50
1:B:619:ILE:HG22	1:B:620:LEU:HD12	1.92	0.50
1:A:347:LEU:O	1:A:348:GLN:HB2	2.12	0.49
1:B:719:ILE:CG2	1:B:723:VAL:HG11	2.34	0.49
1:B:174:MET:HE1	1:B:715:ILE:HG12	1.93	0.49
1:A:616:PHE:CD2	1:A:620:LEU:HD13	2.46	0.49
1:A:685:VAL:O	1:A:689:THR:HG23	2.11	0.49
1:B:86:MET:HG3	1:B:136:PHE:HB3	1.94	0.49
1:A:57:ILE:HG13	1:A:189:PHE:CD2	2.48	0.49
1:B:236:ASP:HB2	1:B:462:PRO:HG3	1.94	0.49
1:B:293:PHE:CE2	1:B:336:THR:HG21	2.46	0.49
1:A:76:THR:HG21	1:A:174:MET:HE2	1.95	0.49
1:A:596:TYR:N	1:A:596:TYR:HD1	2.09	0.49
1:B:327:THR:HG22	1:B:331:LEU:HD12	1.94	0.49
1:A:616:PHE:CE2	1:A:620:LEU:HD13	2.48	0.49
1:B:7:PHE:CA	1:B:10:ILE:HD13	2.41	0.49
1:B:75:GLN:CG	1:B:152:MET:HE1	2.43	0.49
1:A:38:ILE:HD12	1:A:473:MET:HG3	1.95	0.48
1:A:10:ILE:N	1:A:11:PRO:CD	2.76	0.48
1:A:207:LEU:O	1:A:208:VAL:CG1	2.60	0.48
1:B:8:PHE:C	1:B:9:LEU:HD23	2.38	0.48
1:B:39:SER:OG	1:B:103:VAL:HG22	2.13	0.48
1:A:641:THR:HG22	1:A:706:ILE:HG22	1.95	0.48
1:B:199:LYS:HE3	1:B:199:LYS:HA	1.95	0.48
1:A:65:ALA:HB2	1:A:81:LEU:HD13	1.95	0.48
1:A:412:LEU:C	1:A:412:LEU:HD23	2.38	0.48
1:A:245:LEU:HD23	1:A:245:LEU:C	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:VAL:HG22	1:B:650:ILE:HB	1.94	0.48
1:A:86:MET:CE	1:A:139:LEU:HB3	2.44	0.48
1:A:541:ASP:HB3	1:A:544:VAL:HG23	1.95	0.48
1:B:222:ASN:OD1	1:B:222:ASN:O	2.30	0.48
1:B:363:TRP:CD2	1:B:367:ILE:HD11	2.47	0.48
1:A:32:THR:HB	1:A:35:MET:CG	2.44	0.48
1:A:39:SER:O	1:A:42:ILE:HG13	2.14	0.48
1:B:710:SER:O	1:B:714:VAL:HG23	2.14	0.48
1:A:616:PHE:CD1	1:A:616:PHE:C	2.92	0.48
1:A:86:MET:HG3	1:A:136:PHE:HB3	1.95	0.47
1:B:6:LEU:O	1:B:10:ILE:HD12	2.14	0.47
1:A:328:SER:HB2	1:A:453:THR:HG21	1.96	0.47
1:B:10:ILE:N	1:B:11:PRO:CD	2.76	0.47
1:B:401:GLY:O	1:B:405:VAL:HG23	2.14	0.47
1:A:240:LEU:HD13	1:A:458:ASP:CG	2.39	0.47
1:A:723:VAL:O	1:A:723:VAL:CG2	2.62	0.47
1:B:384:THR:HG21	1:B:496:ALA:CB	2.45	0.47
1:A:138:LEU:HD12	1:A:138:LEU:HA	1.66	0.47
1:A:198:THR:HG22	1:A:231:GLY:HA2	1.97	0.47
1:A:262:MET:HE1	1:A:286:LEU:CD1	2.45	0.47
1:A:666:LEU:CD1	1:A:675:GLY:HA2	2.44	0.47
1:B:328:SER:HB2	1:B:453:THR:HG21	1.97	0.47
1:B:613:TYR:HB3	1:B:614:PRO:CD	2.44	0.47
1:B:639:ILE:HD13	1:B:639:ILE:HA	1.62	0.47
1:A:192:VAL:HG11	1:A:614:PRO:HB2	1.97	0.47
1:A:344:LEU:HD22	1:A:347:LEU:HD12	1.96	0.47
1:A:410:LEU:HD13	1:A:659:TRP:HH2	1.80	0.47
1:B:105:VAL:O	1:B:108:ALA:HB3	2.14	0.47
1:A:613:TYR:HB3	1:A:614:PRO:CD	2.44	0.47
1:B:42:ILE:HD12	1:B:42:ILE:C	2.40	0.47
1:B:561:TYR:CD1	1:B:561:TYR:C	2.93	0.47
1:A:225:THR:O	1:A:227:ALA:N	2.48	0.47
1:B:499:LYS:HE2	1:B:695:LYS:O	2.15	0.47
1:A:32:THR:HG22	1:A:34:ARG:N	2.30	0.46
1:A:236:ASP:HB2	1:A:462:PRO:HG3	1.97	0.46
1:A:223:PRO:C	1:A:225:THR:N	2.73	0.46
1:A:263:PHE:HB3	1:A:264:PRO:HD3	1.98	0.46
1:B:386:TYR:HB2	1:B:665:TYR:CD1	2.50	0.46
1:B:240:LEU:HD13	1:B:458:ASP:CB	2.43	0.46
1:A:267:VAL:HG21	1:A:527:ILE:HG23	1.98	0.46
1:A:293:PHE:CE2	1:A:336:THR:HG21	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:LEU:HB3	1:A:347:LEU:HD12	1.98	0.46
1:B:30:GLU:CA	1:B:107:GLU:OE1	2.64	0.46
1:B:191:ARG:CZ	1:B:700:PRO:HB3	2.46	0.46
1:B:192:VAL:HG11	1:B:614:PRO:CB	2.44	0.46
1:B:666:LEU:CD1	1:B:675:GLY:HA2	2.46	0.46
1:A:6:LEU:C	1:A:6:LEU:HD13	2.41	0.46
1:A:198:THR:HG22	1:A:231:GLY:CA	2.46	0.46
1:B:337:ALA:HB2	1:B:363:TRP:NE1	2.30	0.46
1:B:75:GLN:HG2	1:B:152:MET:HE1	1.96	0.46
1:B:692:ASP:OD1	1:B:695:LYS:NZ	2.33	0.46
1:B:86:MET:HE2	1:B:140:GLY:N	2.31	0.46
1:B:138:LEU:HD11	1:B:248:PHE:CZ	2.50	0.45
1:B:288:SER:OG	1:B:343:TYR:OH	2.16	0.45
1:B:57:ILE:HG13	1:B:189:PHE:CD2	2.51	0.45
1:A:121:VAL:CG2	1:A:122:ALA:N	2.79	0.45
1:A:460:TYR:CD1	1:A:460:TYR:C	2.92	0.45
1:B:167:ILE:HG23	1:B:718:SER:HB3	1.99	0.45
1:B:719:ILE:C	1:B:723:VAL:HG12	2.39	0.45
1:B:246:GLU:OE1	1:B:704:ILE:HD12	2.16	0.45
1:A:13:VAL:HG22	1:A:302:MET:SD	2.57	0.45
1:A:445:LEU:O	1:A:449:SER:N	2.50	0.45
1:B:269:LYS:C	1:B:270:ILE:HD12	2.41	0.45
1:B:192:VAL:HG13	1:B:562:LEU:HD21	1.99	0.45
1:B:262:MET:SD	1:B:353:LEU:HD11	2.57	0.45
1:B:527:ILE:N	1:B:527:ILE:CD1	2.80	0.45
1:B:611:MET:O	1:B:612:GLY:C	2.60	0.45
1:A:6:LEU:HD12	1:A:7:PHE:HD1	1.82	0.45
1:A:87:SER:HB2	1:A:186:ILE:HG12	1.99	0.45
1:A:111:THR:HG22	1:A:112:THR:N	2.30	0.45
1:A:191:ARG:CZ	1:A:700:PRO:HB3	2.47	0.45
1:A:267:VAL:HG21	1:A:527:ILE:CG2	2.46	0.45
1:A:511:LEU:HD11	1:B:549:LEU:HD11	1.97	0.45
1:B:573:MET:O	1:B:576:GLU:N	2.50	0.45
1:A:301:SER:HA	1:A:328:SER:OG	2.17	0.45
1:B:75:GLN:OE1	1:B:152:MET:HE1	2.17	0.45
1:B:568:LYS:HD2	1:B:610:GLN:OE1	2.16	0.45
1:B:679:GLU:N	1:B:680:PRO:CD	2.80	0.45
1:B:160:ILE:CD1	1:B:160:ILE:N	2.75	0.44
1:B:573:MET:HE3	1:B:599:CYS:SG	2.57	0.44
1:B:263:PHE:HB3	1:B:264:PRO:HD3	1.97	0.44
1:A:386:TYR:HB2	1:A:665:TYR:CD1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:536:LEU:HD12	1:B:536:LEU:C	2.43	0.44
1:A:627:PHE:HE1	1:A:723:VAL:CG1	2.30	0.44
1:B:25:VAL:HB	1:B:97:MET:HE1	2.00	0.44
1:B:392:GLN:O	1:B:396:LYS:HG3	2.18	0.44
1:B:666:LEU:HD13	1:B:675:GLY:HA2	2.00	0.44
1:A:170:VAL:HG23	1:A:261:TYR:CE1	2.53	0.44
1:B:222:ASN:N	1:B:223:PRO:CD	2.81	0.44
1:B:262:MET:SD	1:B:353:LEU:CD1	3.06	0.44
1:B:185:ILE:C	1:B:185:ILE:CD1	2.80	0.44
1:A:425:VAL:HG13	1:B:546:ALA:HB1	2.00	0.43
1:B:42:ILE:HD11	1:B:102:ASN:HD21	1.82	0.43
1:B:87:SER:HB2	1:B:186:ILE:HG12	1.99	0.43
1:B:256:ILE:HG23	1:B:287:ILE:HG23	1.99	0.43
1:B:675:GLY:O	1:B:676:LYS:HB3	2.17	0.43
1:A:174:MET:HE3	1:A:174:MET:HA	2.00	0.43
1:A:174:MET:HE3	1:A:174:MET:O	2.17	0.43
1:A:332:THR:O	1:A:336:THR:HB	2.19	0.43
1:A:199:LYS:HG2	1:A:697:THR:OG1	2.19	0.43
1:B:579:ARG:O	1:B:583:GLU:HB2	2.18	0.43
1:B:608:LEU:HD23	1:B:611:MET:CE	2.49	0.43
1:B:725:LEU:HD12	1:B:726:PHE:CE2	2.54	0.43
1:A:118:ALA:C	1:A:121:VAL:HG22	2.41	0.43
1:B:13:VAL:HG22	1:B:302:MET:SD	2.59	0.43
1:B:301:SER:HA	1:B:328:SER:OG	2.18	0.43
1:B:263:PHE:HB3	1:B:264:PRO:CD	2.49	0.43
1:B:347:LEU:O	1:B:348:GLN:HB2	2.19	0.43
1:A:499:LYS:HE2	1:A:695:LYS:O	2.19	0.43
1:B:484:THR:HG22	1:B:485:ASP:N	2.33	0.43
1:A:679:GLU:N	1:A:680:PRO:CD	2.82	0.42
1:B:12:LEU:CD1	1:B:139:LEU:HD13	2.49	0.42
1:A:240:LEU:HD13	1:A:458:ASP:OD2	2.19	0.42
1:B:363:TRP:CH2	1:B:367:ILE:CD1	3.02	0.42
1:B:42:ILE:CD1	1:B:102:ASN:HD21	2.32	0.42
1:B:198:THR:HG22	1:B:199:LYS:NZ	2.34	0.42
1:B:208:VAL:HG13	1:B:208:VAL:O	2.19	0.42
1:B:332:THR:O	1:B:336:THR:HB	2.18	0.42
1:B:507:ILE:HD13	1:B:646:ALA:HB1	2.01	0.42
1:A:38:ILE:CD1	1:A:470:ILE:HA	2.50	0.42
1:A:523:SER:O	1:A:524:PRO:C	2.61	0.42
1:B:7:PHE:O	1:B:10:ILE:HD13	2.20	0.42
1:A:6:LEU:C	1:A:6:LEU:CD1	2.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:ILE:CG2	4:B:2004:HOH:O	2.57	0.42
1:B:666:LEU:C	1:B:666:LEU:CD1	2.91	0.42
1:A:129:MET:O	1:A:133:VAL:HG23	2.19	0.42
1:A:199:LYS:HA	1:A:199:LYS:HE3	2.02	0.42
1:B:389:LYS:N	1:B:390:PRO:CD	2.82	0.42
1:B:499:LYS:NZ	1:B:695:LYS:O	2.52	0.42
1:A:207:LEU:C	1:A:208:VAL:HG13	2.45	0.42
1:A:347:LEU:HA	1:A:347:LEU:HD23	1.81	0.42
1:B:262:MET:CE	1:B:353:LEU:HD11	2.50	0.42
1:A:65:ALA:HB2	1:A:81:LEU:CD1	2.50	0.42
1:A:158:LEU:HD21	1:A:283:ILE:HG21	2.01	0.42
1:A:263:PHE:HB3	1:A:264:PRO:CD	2.50	0.42
1:A:87:SER:CB	1:A:186:ILE:HG12	2.50	0.42
1:A:100:ARG:O	1:A:103:VAL:HG23	2.19	0.42
1:A:344:LEU:CD2	1:A:347:LEU:HD12	2.50	0.42
1:A:417:VAL:HG22	1:A:650:ILE:HB	2.02	0.42
1:B:115:ILE:HA	1:B:483:ILE:HG21	2.02	0.42
1:B:582:ARG:HD2	1:B:582:ARG:HA	1.90	0.42
1:A:39:SER:HB2	1:A:103:VAL:CG1	2.48	0.41
1:A:401:GLY:O	1:A:405:VAL:HG23	2.20	0.41
1:A:666:LEU:C	1:A:666:LEU:CD1	2.93	0.41
1:B:674:TYR:O	1:B:675:GLY:C	2.60	0.41
1:B:39:SER:O	1:B:42:ILE:HG13	2.19	0.41
1:B:725:LEU:CD1	1:B:726:PHE:CD1	3.03	0.41
1:B:82:LEU:O	1:B:86:MET:HB2	2.20	0.41
1:B:507:ILE:CD1	1:B:646:ALA:HB1	2.51	0.41
1:A:167:ILE:HG23	1:A:718:SER:HB3	2.02	0.41
1:B:147:ILE:O	1:B:151:TRP:HB3	2.21	0.41
1:A:344:LEU:CD2	1:A:347:LEU:HD11	2.51	0.41
1:A:573:MET:O	1:A:574:VAL:C	2.63	0.41
1:B:75:GLN:H	1:B:75:GLN:NE2	2.11	0.41
1:A:192:VAL:HG11	1:A:614:PRO:CB	2.50	0.41
1:A:448:LEU:CD2	1:A:505:SER:HB3	2.51	0.41
1:B:75:GLN:HG2	1:B:152:MET:CE	2.51	0.41
1:A:82:LEU:HD23	1:A:82:LEU:HA	1.92	0.41
1:A:513:LEU:HD23	1:A:513:LEU:HA	1.94	0.41
1:A:639:ILE:HD13	1:A:639:ILE:HA	1.78	0.41
1:B:87:SER:CB	1:B:186:ILE:HG12	2.51	0.41
1:B:6:LEU:O	1:B:10:ILE:CD1	2.70	0.40
1:A:333:VAL:HG13	1:A:367:ILE:HD11	2.03	0.40
1:A:606:ASN:O	1:A:606:ASN:ND2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:674:TYR:CD1	1:B:680:PRO:HG2	2.56	0.40
1:A:639:ILE:HD12	1:A:639:ILE:HG23	1.95	0.40
1:B:87:SER:O	1:B:88:ALA:C	2.64	0.40
1:B:138:LEU:CD1	1:B:248:PHE:CZ	3.05	0.40
1:B:221:ARG:C	1:B:223:PRO:HD3	2.45	0.40
1:B:328:SER:OG	1:B:453:THR:HG21	2.20	0.40
1:A:344:LEU:HA	1:A:347:LEU:CD1	2.52	0.40
1:B:448:LEU:CD2	1:B:505:SER:HB3	2.51	0.40
1:B:336:THR:CG2	1:B:363:TRP:CD1	2.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	686/735 (93%)	632 (92%)	45 (7%)	9 (1%)	9	21
1	B	699/735 (95%)	653 (93%)	33 (5%)	13 (2%)	6	13
All	All	1385/1470 (94%)	1285 (93%)	78 (6%)	22 (2%)	7	16

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	359	ALA
1	B	32	THR
1	B	359	ALA
1	B	675	GLY
1	B	676	LYS
1	A	208	VAL
1	A	226	ILE
1	A	387	ARG
1	A	450	PHE

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Mol	Chain	Res	Type
1	A	672	GLU
1	A	678	SER
1	B	223	PRO
1	B	358	GLY
1	B	387	ARG
1	A	477	ASP
1	A	478	PRO
1	B	477	ASP
1	B	478	PRO
1	B	589	GLU
1	B	584	ILE
1	B	585	PRO
1	B	600	ILE

5.3.2 Protein sidechains ⓘ

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	470/575 (82%)	406 (86%)	64 (14%)	3	7
1	B	486/575 (84%)	425 (87%)	61 (13%)	4	9
All	All	956/1150 (83%)	831 (87%)	125 (13%)	4	8

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	13	VAL
1	A	38	ILE
1	A	50	LEU
1	A	54	THR
1	A	67	LEU
1	A	81	LEU
1	A	82	LEU
1	A	86	MET
1	A	103	VAL

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Mol	Chain	Res	Type
1	A	105	VAL
1	A	110	ARG
1	A	111	THR
1	A	112	THR
1	A	138	LEU
1	A	158	LEU
1	A	160	ILE
1	A	162	THR
1	A	186	ILE
1	A	199	LYS
1	A	225	THR
1	A	240	LEU
1	A	262	MET
1	A	267	VAL
1	A	270	ILE
1	A	274	LEU
1	A	298	LEU
1	A	350	LEU
1	A	367	ILE
1	A	384	THR
1	A	407	SER
1	A	410	LEU
1	A	417	VAL
1	A	422	LEU
1	A	429	LEU
1	A	437	LEU
1	A	448	LEU
1	A	459	SER
1	A	473	MET
1	A	484	THR
1	A	490	VAL
1	A	508	PHE
1	A	522	ILE
1	A	537	LEU
1	A	540	LEU
1	A	550	LEU
1	A	554	ILE
1	A	564	SER
1	A	571	MET
1	A	596	TYR
1	A	599	CYS
1	A	603	THR

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Mol	Chain	Res	Type
1	A	613	TYR
1	A	638	LEU
1	A	639	ILE
1	A	643	LEU
1	A	664	LYS
1	A	666	LEU
1	A	679	GLU
1	A	684	LEU
1	A	689	THR
1	A	702	LEU
1	A	705	LEU
1	A	719	ILE
1	B	9	LEU
1	B	13	VAL
1	B	54	THR
1	B	75	GLN
1	B	81	LEU
1	B	82	LEU
1	B	86	MET
1	B	103	VAL
1	B	110	ARG
1	B	111	THR
1	B	138	LEU
1	B	143	LEU
1	B	160	ILE
1	B	162	THR
1	B	185	ILE
1	B	186	ILE
1	B	198	THR
1	B	199	LYS
1	B	203	MET
1	B	225	THR
1	B	230	VAL
1	B	240	LEU
1	B	262	MET
1	B	267	VAL
1	B	298	LEU
1	B	318	GLN
1	B	331	LEU
1	B	336	THR
1	B	347	LEU
1	B	353	LEU

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Mol	Chain	Res	Type
1	B	384	THR
1	B	407	SER
1	B	410	LEU
1	B	422	LEU
1	B	429	LEU
1	B	437	LEU
1	B	448	LEU
1	B	470	ILE
1	B	484	THR
1	B	508	PHE
1	B	522	ILE
1	B	527	ILE
1	B	537	LEU
1	B	540	LEU
1	B	550	LEU
1	B	554	ILE
1	B	564	SER
1	B	571	MET
1	B	583	GLU
1	B	588	LEU
1	B	638	LEU
1	B	639	ILE
1	B	643	LEU
1	B	655	SER
1	B	666	LEU
1	B	672	GLU
1	B	679	GLU
1	B	684	LEU
1	B	689	THR
1	B	702	LEU
1	B	705	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	GLN
1	A	276	HIS
1	A	392	GLN
1	A	606	ASN
1	B	75	GLN
1	B	222	ASN
1	B	318	GLN

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Mol	Chain	Res	Type
1	B	392	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	694/735 (94%)	-0.03	17 (2%) 59 54	48, 84, 143, 186	0
1	B	707/735 (96%)	-0.03	16 (2%) 61 55	49, 87, 145, 219	0
All	All	1401/1470 (95%)	-0.03	33 (2%) 59 54	48, 86, 145, 219	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	TYR	4.8
1	A	391	THR	4.4
1	A	453	THR	3.9
1	B	382	TYR	3.7
1	A	395	GLY	3.3
1	A	224	ALA	3.1
1	A	452	ALA	3.1
1	A	40	SER	3.1
1	A	38	ILE	3.0
1	B	672	GLU	2.9
1	A	392	GLN	2.9
1	B	91	GLY	2.8
1	A	116	GLY	2.7
1	B	590	GLY	2.6
1	A	480	VAL	2.6
1	B	204	ALA	2.6
1	A	599	CYS	2.5
1	B	599	CYS	2.5
1	B	199	LYS	2.5
1	B	115	ILE	2.5
1	B	206	ASP	2.4
1	A	27	ARG	2.3
1	B	156	ASP	2.3
1	A	91	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	228	ASP	2.1
1	B	386	TYR	2.1
1	B	192	VAL	2.1
1	A	342	PHE	2.1
1	B	495	ALA	2.1
1	A	719	ILE	2.1
1	B	532	SER	2.0
1	A	381	GLU	2.0
1	B	205	ALA	2.0

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

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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	727	1/1	0.80	0.13	105,105,105,105	0
2	MG	B	727	1/1	0.80	0.09	96,96,96,96	0
3	CA	B	728	1/1	0.88	0.16	98,98,98,98	0
3	CA	A	728	1/1	0.97	0.04	107,107,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.