



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:55 PM UTC

PDB ID : 5YZV / pdb_00005yzv
Title : Biophysical and structural characterization of the thermostable WD40 domain of a prokaryotic protein, Thermomonospora curvata PkwA
Authors : Li, D.Y.; Shen, C.; Du, Y.; Qiao, F.F.; Kong, T.; Yuan, L.R.; Zhang, D.L.; Wu, X.H.; Wu, Y.D.
Deposited on : 2017-12-15
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

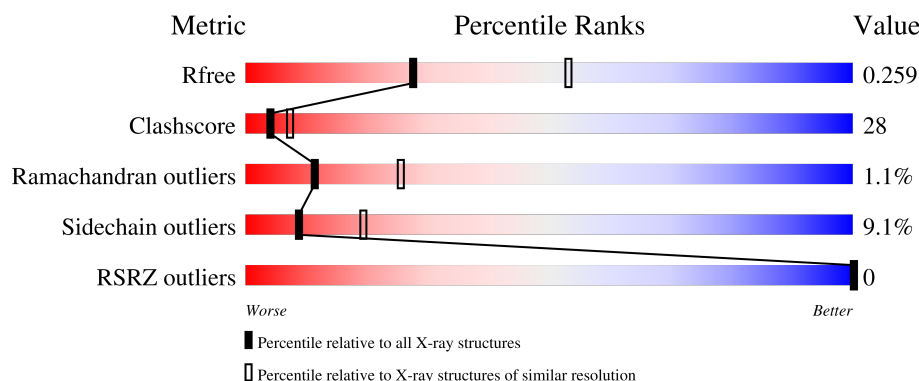
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

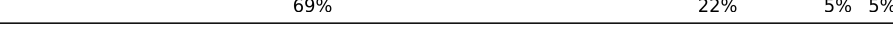
The reported resolution of this entry is 2.60 Å.

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	302	
1	B	302	
1	C	302	
1	D	302	
1	E	302	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10922 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 Probable serine/threonine-protein kinase PkwA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2124	1325	369	427	3			
1	B	288	Total	C	N	O	S	0	0	0
			2154	1341	377	433	3			
1	C	290	Total	C	N	O	S	0	0	0
			2152	1341	373	435	3			
1	D	284	Total	C	N	O	S	0	0	0
			2111	1319	364	425	3			
1	E	288	Total	C	N	O	S	0	0	0
			2151	1340	376	432	3			

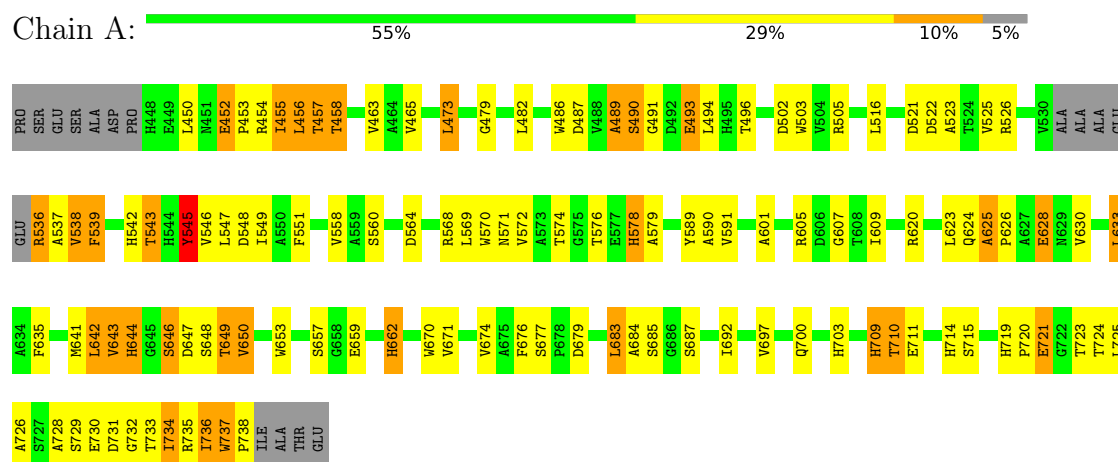
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	57	Total	O	0	0
			57	57		
2	B	44	Total	O	0	0
			44	44		
2	C	44	Total	O	0	0
			44	44		
2	D	42	Total	O	0	0
			42	42		
2	E	43	Total	O	0	0
			43	43		

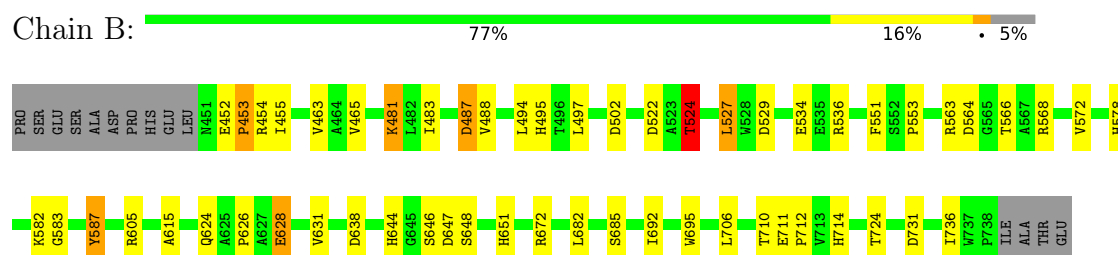
3 Residue-property plots

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.

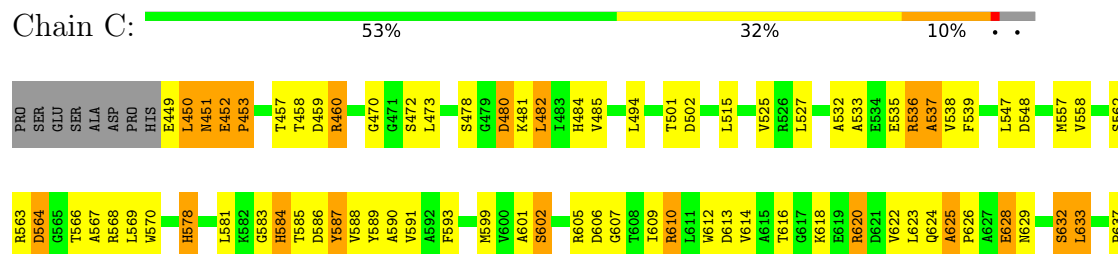
- Molecule 1: Probable serine/threonine-protein kinase PkwA

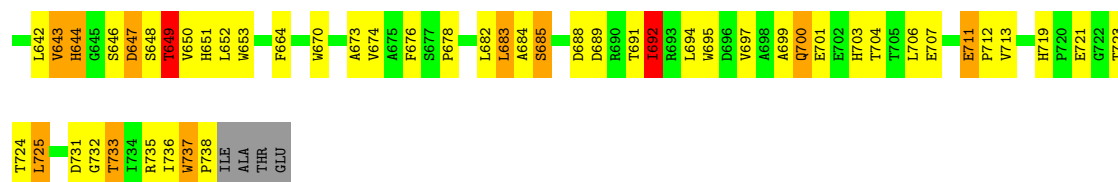


- Molecule 1: Probable serine/threonine-protein kinase PkwA



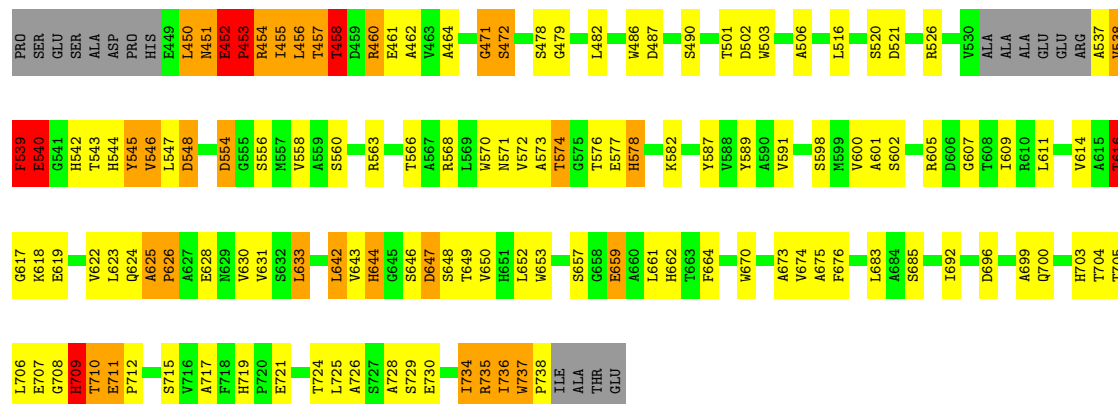
- Molecule 1: Probable serine/threonine-protein kinase PkwA





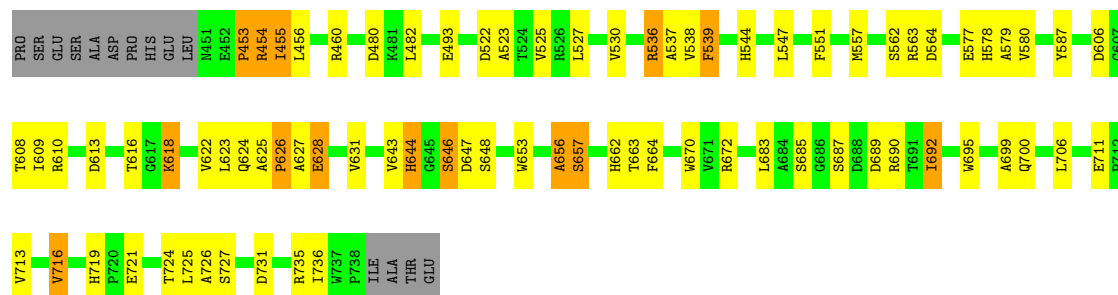
- Molecule 1: Probable serine/threonine-protein kinase PkwA

Chain D: 50% 32% 10% 6%



- Molecule 1: Probable serine/threonine-protein kinase PkwA

Chain E: 69% 22% 5% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	43.69Å 107.26Å 110.27Å 78.77° 89.33° 88.91°	Depositor
Resolution (Å)	50.01 – 2.60 50.01 – 2.60	Depositor EDS
% Data completeness (in resolution range)	90.1 (50.01-2.60) 90.1 (50.01-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.62 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.225 , 0.259 0.230 , 0.259	Depositor DCC
R_{free} test set	2895 reflections (4.28%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 24.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l 0.013 for -h,l,k 0.118 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10922	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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	1.46	17/2178 (0.8%)	1.31	20/2981 (0.7%)
1	B	1.29	7/2209 (0.3%)	1.13	8/3022 (0.3%)
1	C	1.45	20/2207 (0.9%)	1.39	17/3022 (0.6%)
1	D	1.49	22/2165 (1.0%)	1.33	19/2964 (0.6%)
1	E	1.29	15/2206 (0.7%)	1.05	5/3018 (0.2%)
All	All	1.40	81/10965 (0.7%)	1.25	69/15007 (0.5%)

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	453	PRO	N-CA	16.78	1.69	1.47
1	D	539	PHE	CA-C	-12.39	1.37	1.52
1	D	455	ILE	CA-CB	-10.52	1.42	1.54
1	A	538	VAL	CA-CB	-9.28	1.42	1.54
1	A	455	ILE	CA-CB	-9.15	1.43	1.54
1	C	452	GLU	C-N	9.15	1.45	1.34
1	A	456	LEU	CA-C	-8.60	1.42	1.52
1	C	649	THR	CB-CG2	-8.34	1.25	1.52
1	A	491	GLY	N-CA	-7.72	1.35	1.45
1	A	536	ARG	CA-C	-7.53	1.37	1.52
1	A	538	VAL	N-CA	-7.28	1.38	1.46
1	E	454	ARG	CA-CB	-7.17	1.42	1.53
1	D	454	ARG	N-CA	-7.12	1.37	1.46
1	C	533	ALA	N-CA	-7.11	1.36	1.46
1	B	524	THR	CB-CG2	-6.93	1.29	1.52
1	C	535	GLU	CA-C	-6.88	1.43	1.52
1	D	455	ILE	N-CA	-6.81	1.38	1.46
1	C	537	ALA	N-CA	-6.78	1.36	1.45
1	D	454	ARG	CA-C	-6.67	1.44	1.52
1	A	455	ILE	N-CA	-6.63	1.38	1.46
1	C	628	GLU	C-O	-6.63	1.17	1.24
1	E	626	PRO	N-CA	-6.56	1.39	1.46
1	B	453	PRO	C-O	-6.55	1.15	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	453	PRO	CA-C	-6.53	1.43	1.52
1	E	626	PRO	CA-CB	-6.53	1.45	1.53
1	A	490	SER	CA-C	-6.49	1.43	1.52
1	D	453	PRO	N-CA	-6.45	1.39	1.47
1	E	538	VAL	N-CA	-6.44	1.38	1.46
1	D	538	VAL	CA-CB	-6.31	1.45	1.54
1	C	584	HIS	N-CA	-6.31	1.39	1.46
1	D	452	GLU	CA-C	-6.26	1.45	1.52
1	C	532	ALA	CA-C	-6.25	1.45	1.52
1	D	538	VAL	C-O	-6.22	1.15	1.24
1	D	452	GLU	CA-CB	-6.15	1.43	1.53
1	D	540	GLU	CA-C	-6.10	1.45	1.52
1	D	539	PHE	CB-CG	-6.03	1.36	1.50
1	C	692	ILE	CA-C	-5.97	1.45	1.52
1	A	456	LEU	N-CA	-5.96	1.38	1.45
1	B	455	ILE	N-CA	-5.95	1.39	1.46
1	E	455	ILE	N-CA	-5.89	1.39	1.46
1	C	593	PHE	C-O	-5.82	1.16	1.23
1	A	649	THR	CA-CB	-5.79	1.43	1.53
1	D	574	THR	CB-CG2	-5.74	1.33	1.52
1	A	709	HIS	C-O	-5.71	1.16	1.23
1	C	538	VAL	N-CA	-5.63	1.39	1.46
1	B	536	ARG	CA-CB	-5.63	1.45	1.53
1	E	656	ALA	C-O	-5.58	1.17	1.23
1	D	460	ARG	C-O	-5.56	1.19	1.24
1	A	574	THR	CB-CG2	-5.51	1.34	1.52
1	D	616	THR	C-O	-5.49	1.17	1.24
1	A	493	GLU	CA-C	-5.44	1.46	1.52
1	E	454	ARG	N-CA	-5.44	1.39	1.45
1	B	626	PRO	CA-C	-5.39	1.44	1.52
1	E	455	ILE	CA-C	-5.36	1.46	1.52
1	C	451	ASN	C-O	-5.33	1.17	1.23
1	E	455	ILE	C-O	-5.32	1.18	1.24
1	D	548	ASP	CB-CG	-5.30	1.38	1.52
1	C	460	ARG	CG-CD	-5.30	1.36	1.52
1	D	455	ILE	C-O	-5.29	1.18	1.24
1	A	576	THR	CB-CG2	-5.28	1.35	1.52
1	E	580	VAL	C-O	-5.28	1.18	1.24
1	C	538	VAL	C-O	-5.26	1.18	1.24
1	A	538	VAL	CA-C	-5.25	1.46	1.52
1	C	733	THR	CB-CG2	-5.25	1.35	1.52
1	E	537	ALA	CA-CB	-5.25	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	454	ARG	CA-C	-5.23	1.46	1.52
1	E	456	LEU	CA-C	-5.20	1.46	1.52
1	D	464	ALA	N-CA	-5.19	1.39	1.46
1	C	700	GLN	CA-C	5.17	1.59	1.52
1	D	709	HIS	CA-C	-5.16	1.45	1.52
1	C	647	ASP	CB-CG	5.13	1.64	1.52
1	E	539	PHE	CA-CB	-5.11	1.45	1.53
1	B	454	ARG	CA-CB	-5.10	1.45	1.53
1	C	480	ASP	C-O	-5.08	1.17	1.23
1	A	578	HIS	CA-C	-5.06	1.45	1.52
1	D	626	PRO	CA-CB	-5.05	1.47	1.53
1	E	536	ARG	N-CA	-5.03	1.39	1.46
1	C	536	ARG	N-CA	-5.02	1.38	1.46
1	A	456	LEU	CA-CB	-5.02	1.43	1.54
1	D	458	THR	CB-CG2	-5.02	1.35	1.52
1	D	460	ARG	CA-CB	-5.00	1.47	1.54

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	452	GLU	CA-C-N	19.01	139.57	118.85
1	C	452	GLU	C-N-CA	19.01	139.57	118.85
1	C	453	PRO	CA-N-CD	-10.64	97.10	112.00
1	D	455	ILE	N-CA-C	9.19	121.04	108.17
1	D	456	LEU	N-CA-C	8.85	123.48	109.50
1	A	650	VAL	N-CA-C	8.71	120.19	109.30
1	A	489	ALA	N-CA-C	-8.52	98.49	110.50
1	C	564	ASP	CA-CB-CG	8.26	120.86	112.60
1	C	628	GLU	N-CA-C	8.03	118.84	108.24
1	A	455	ILE	N-CA-C	7.90	119.53	107.99
1	C	610	ARG	NE-CZ-NH1	-7.82	113.68	121.50
1	B	553	PRO	CA-C-N	7.40	130.19	120.28
1	B	553	PRO	C-N-CA	7.40	130.19	120.28
1	C	480	ASP	N-CA-C	-7.27	100.94	110.53
1	D	471	GLY	CA-C-N	7.07	134.42	121.70
1	D	471	GLY	C-N-CA	7.07	134.42	121.70
1	C	620	ARG	N-CA-C	6.89	118.87	111.36
1	D	538	VAL	CB-CA-C	-6.77	101.52	111.19
1	A	455	ILE	N-CA-CB	-6.71	103.36	111.21
1	C	449	GLU	CA-C-N	6.66	134.26	121.54
1	C	449	GLU	C-N-CA	6.66	134.26	121.54
1	A	538	VAL	CB-CA-C	-6.49	102.01	111.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	709	HIS	N-CA-C	6.24	124.08	110.80
1	A	578	HIS	N-CA-C	6.23	121.01	113.41
1	A	543	THR	N-CA-C	-6.19	104.59	111.71
1	D	710	THR	N-CA-C	-6.16	104.68	111.82
1	C	610	ARG	NE-CZ-NH2	6.10	124.69	119.20
1	E	628	GLU	N-CA-C	6.07	118.52	108.99
1	D	455	ILE	N-CA-CB	-6.01	103.08	111.25
1	C	537	ALA	N-CA-C	5.84	117.87	110.33
1	A	649	THR	N-CA-C	5.72	117.23	108.42
1	A	578	HIS	CB-CA-C	-5.72	99.23	109.24
1	D	452	GLU	N-CA-C	5.70	122.40	109.81
1	A	646	SER	N-CA-C	5.66	116.83	108.86
1	D	460	ARG	NE-CZ-NH1	-5.61	115.89	121.50
1	D	649	THR	N-CA-C	5.57	117.09	109.18
1	A	456	LEU	N-CA-C	5.55	118.13	109.52
1	C	564	ASP	CB-CA-C	-5.51	99.64	110.11
1	D	539	PHE	CA-CB-CG	-5.47	108.33	113.80
1	C	647	ASP	CB-CA-C	5.45	120.41	111.41
1	D	453	PRO	CB-CA-C	-5.41	102.63	111.56
1	E	625	ALA	CA-C-O	5.38	127.47	121.22
1	A	536	ARG	CB-CA-C	-5.38	99.88	110.10
1	B	536	ARG	N-CA-CB	-5.37	103.44	110.59
1	D	673	ALA	N-CA-C	5.37	117.55	109.23
1	B	605	ARG	NE-CZ-NH2	5.35	124.02	119.20
1	E	687	SER	N-CA-C	5.35	117.54	109.41
1	D	460	ARG	N-CA-C	5.29	115.01	108.19
1	B	452	GLU	CA-C-N	-5.26	113.26	119.84
1	B	452	GLU	C-N-CA	-5.26	113.26	119.84
1	D	539	PHE	CB-CA-C	-5.23	97.95	109.56
1	A	643	VAL	N-CA-C	5.21	115.40	108.11
1	A	538	VAL	CA-CB-CG2	-5.21	101.55	110.40
1	D	478	SER	N-CA-C	5.19	117.30	109.41
1	A	662	HIS	N-CA-C	5.17	116.15	108.60
1	A	687	SER	N-CA-C	5.16	117.03	109.24
1	A	454	ARG	CB-CA-C	5.14	118.52	110.19
1	E	453	PRO	CB-CA-C	-5.14	103.08	111.56
1	A	545	TYR	CA-C-N	-5.14	116.82	123.19
1	A	545	TYR	C-N-CA	-5.14	116.82	123.19
1	C	707	GLU	N-CA-C	5.13	119.71	111.81
1	E	530	VAL	CB-CA-C	-5.13	105.30	112.02
1	B	488	VAL	N-CA-C	5.13	115.34	110.42
1	D	453	PRO	CA-N-CD	-5.09	104.87	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	700	GLN	N-CA-CB	-5.08	104.20	113.10
1	D	647	ASP	N-CA-C	5.03	118.52	112.38
1	C	578	HIS	CA-CB-CG	-5.02	108.78	113.80
1	B	487	ASP	N-CA-C	-5.02	101.57	109.25
1	A	649	THR	CB-CA-C	-5.00	102.03	110.43

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2124	0	1969	154	0
1	B	2154	0	2013	30	0
1	C	2152	0	1995	170	0
1	D	2111	0	1963	174	0
1	E	2151	0	2009	48	0
2	A	57	0	0	3	0
2	B	44	0	0	0	0
2	C	44	0	0	4	0
2	D	42	0	0	0	0
2	E	43	0	0	0	0
All	All	10922	0	9949	570	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 28.

All (570) 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:452:GLU:HG2	1:A:453:PRO:CD	1.31	1.57
1:C:453:PRO:N	1:C:453:PRO:CA	1.69	1.34
1:A:452:GLU:CG	1:A:453:PRO:CD	2.13	1.27
1:A:452:GLU:CG	1:A:453:PRO:HD3	1.64	1.24
1:C:599:MET:HE1	1:C:620:ARG:CD	1.71	1.19

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:453:PRO:C	1:D:454:ARG:HD2	1.70	1.16
1:C:605:ARG:HG3	1:C:629:ASN:ND2	1.60	1.16
1:C:694:LEU:HD21	1:C:725:LEU:CD2	1.77	1.15
1:D:662:HIS:HB3	1:D:700:GLN:CD	1.71	1.14
1:C:599:MET:HE1	1:C:620:ARG:HD2	1.31	1.12
1:D:662:HIS:HB3	1:D:700:GLN:OE1	1.49	1.10
1:C:643:VAL:HG21	1:C:683:LEU:HD11	1.35	1.06
1:A:456:LEU:O	1:A:457:THR:HG23	1.55	1.05
1:C:694:LEU:HD21	1:C:725:LEU:HD21	1.32	1.04
1:A:643:VAL:HG21	1:A:683:LEU:HD11	1.39	1.02
1:C:605:ARG:HG3	1:C:629:ASN:HD21	1.25	1.02
1:A:452:GLU:CG	1:A:453:PRO:HD2	1.94	0.98
1:B:487:ASP:HB2	1:B:494:LEU:HD11	1.43	0.97
1:C:599:MET:HE1	1:C:620:ARG:HD3	1.46	0.95
1:D:571:ASN:HB2	1:D:578:HIS:HE1	1.31	0.94
1:C:628:GLU:HG2	1:C:629:ASN:H	1.33	0.93
1:C:451:ASN:O	1:C:738:PRO:HD2	1.69	0.90
1:D:650:VAL:HG12	1:D:664:PHE:HB2	1.54	0.90
1:C:643:VAL:HG21	1:C:683:LEU:CD1	2.02	0.89
1:D:643:VAL:HG12	1:D:652:LEU:HD23	1.55	0.88
1:C:599:MET:CE	1:C:620:ARG:HD3	2.03	0.88
1:C:451:ASN:O	1:C:738:PRO:CD	2.22	0.87
1:C:692:ILE:CG1	1:C:706:LEU:HB2	2.04	0.87
1:A:452:GLU:CB	1:A:453:PRO:HD2	2.04	0.87
1:C:450:LEU:CB	1:C:704:THR:HG21	2.05	0.86
1:D:461:GLU:HG3	1:D:503:TRP:HH2	1.40	0.85
1:A:607:GLY:O	1:A:626:PRO:HG2	1.77	0.84
1:C:599:MET:CE	1:C:620:ARG:CD	2.54	0.84
1:D:453:PRO:CD	1:D:736:ILE:HG13	2.06	0.84
1:D:650:VAL:CG1	1:D:664:PHE:HB2	2.10	0.81
1:A:452:GLU:HB3	1:A:453:PRO:HD2	1.63	0.81
1:C:460:ARG:HD2	1:C:478:SER:OG	1.81	0.81
1:A:643:VAL:HG11	1:A:683:LEU:HD21	1.61	0.80
1:D:542:HIS:CE1	1:D:568:ARG:HH21	1.99	0.80
1:A:643:VAL:HG21	1:A:683:LEU:CD1	2.10	0.80
1:D:451:ASN:ND2	1:D:451:ASN:O	2.14	0.80
1:A:456:LEU:O	1:A:457:THR:CG2	2.28	0.80
1:D:571:ASN:HB2	1:D:578:HIS:CE1	2.16	0.80
1:C:616:THR:HG23	1:C:618:LYS:HB2	1.62	0.79
1:D:460:ARG:HG2	1:D:461:GLU:H	1.47	0.79
1:B:615:ALA:HB1	1:C:719:HIS:HE2	1.47	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:482:LEU:HD13	1:C:484:HIS:NE2	1.98	0.79
1:C:699:ALA:O	1:C:700:GLN:HB3	1.82	0.79
1:C:711:GLU:HG2	1:C:712:PRO:HD2	1.64	0.79
1:D:699:ALA:O	1:D:700:GLN:HB2	1.82	0.79
1:D:708:GLY:O	1:D:709:HIS:HB2	1.82	0.78
1:C:607:GLY:O	1:C:626:PRO:HG2	1.84	0.77
1:D:454:ARG:HD2	1:D:454:ARG:N	1.93	0.77
1:D:453:PRO:HD2	1:D:736:ILE:HG13	1.65	0.77
1:A:452:GLU:CB	1:A:453:PRO:CD	2.58	0.77
1:A:490:SER:O	1:A:490:SER:OG	1.98	0.77
1:A:626:PRO:HB3	1:A:644:HIS:NE2	1.99	0.77
1:E:455:ILE:N	1:E:455:ILE:HD12	2.00	0.76
1:C:723:THR:HG23	1:C:724:THR:HG23	1.67	0.76
1:A:526:ARG:NH1	1:A:539:PHE:HE2	1.83	0.76
1:A:643:VAL:HG23	1:A:676:PHE:CE2	2.21	0.76
1:A:450:LEU:O	1:A:738:PRO:HD3	1.86	0.76
1:D:715:SER:OG	1:D:728:ALA:HB3	1.84	0.76
1:D:539:PHE:CD1	1:D:539:PHE:N	2.49	0.75
1:A:625:ALA:N	1:A:626:PRO:HD2	2.01	0.75
1:D:643:VAL:HG13	1:D:676:PHE:CZ	2.20	0.75
1:D:643:VAL:HG12	1:D:652:LEU:CD2	2.17	0.74
1:C:587:TYR:CE1	1:C:605:ARG:HD3	2.23	0.73
1:C:664:PHE:CD1	1:C:700:GLN:CG	2.72	0.73
1:C:609:ILE:HD11	1:C:644:HIS:HB2	1.70	0.73
1:C:548:ASP:OD2	1:C:591:VAL:N	2.20	0.73
1:D:566:THR:OG1	1:D:568:ARG:NH1	2.22	0.73
1:D:538:VAL:HG12	1:D:539:PHE:N	2.02	0.72
1:C:450:LEU:CB	1:C:704:THR:CG2	2.68	0.72
1:A:538:VAL:HG12	1:A:539:PHE:N	2.04	0.72
1:A:571:ASN:HB2	1:A:578:HIS:NE2	2.04	0.72
1:D:643:VAL:HG13	1:D:676:PHE:CE2	2.23	0.72
1:D:457:THR:HG22	1:D:486:TRP:HH2	1.54	0.72
1:D:650:VAL:HG13	1:D:664:PHE:CD2	2.24	0.72
1:C:570:TRP:HA	1:C:578:HIS:HB3	1.70	0.72
1:C:664:PHE:CD1	1:C:700:GLN:HG2	2.26	0.71
1:A:526:ARG:NH1	1:A:539:PHE:CE2	2.58	0.71
1:C:711:GLU:CG	1:C:712:PRO:HD2	2.20	0.71
1:A:643:VAL:HG23	1:A:676:PHE:CZ	2.25	0.71
1:D:454:ARG:N	1:D:454:ARG:CD	2.51	0.71
1:E:644:HIS:HD2	1:E:653:TRP:HE1	1.39	0.71
1:D:662:HIS:CB	1:D:700:GLN:CD	2.59	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:711:GLU:HB2	1:D:729:SER:OG	1.90	0.70
1:D:607:GLY:O	1:D:626:PRO:HG2	1.91	0.70
1:A:568:ARG:HG2	1:A:568:ARG:HH21	1.55	0.70
1:A:564:ASP:OD2	1:A:568:ARG:NH1	2.20	0.70
1:B:564:ASP:OD1	1:B:566:THR:HG22	1.92	0.70
1:A:542:HIS:CE1	1:A:568:ARG:HG3	2.26	0.70
1:D:626:PRO:HB3	1:D:644:HIS:NE2	2.06	0.69
1:D:460:ARG:CZ	1:D:503:TRP:HZ2	2.04	0.69
1:D:454:ARG:HG3	1:D:735:ARG:HG3	1.74	0.69
1:D:538:VAL:CG1	1:D:539:PHE:N	2.56	0.69
1:D:644:HIS:HB3	1:D:653:TRP:HE1	1.57	0.69
1:A:452:GLU:HG2	1:A:453:PRO:HD3	0.70	0.69
1:C:626:PRO:HB3	1:C:644:HIS:NE2	2.07	0.69
1:D:457:THR:CG2	1:D:486:TRP:HH2	2.06	0.69
1:D:625:ALA:N	1:D:626:PRO:HD2	2.07	0.68
1:A:455:ILE:HB	1:A:734:ILE:HB	1.75	0.68
1:A:643:VAL:CG1	1:A:683:LEU:HD21	2.23	0.68
1:B:487:ASP:CB	1:B:494:LEU:HD11	2.23	0.68
1:C:453:PRO:HD3	1:C:738:PRO:HD3	1.75	0.68
1:A:648:SER:O	1:A:649:THR:HG23	1.93	0.68
1:C:673:ALA:HB1	2:C:834:HOH:O	1.93	0.68
1:D:719:HIS:ND1	1:D:721:GLU:HG2	2.09	0.68
1:A:719:HIS:CE1	1:A:721:GLU:HG3	2.29	0.67
1:D:452:GLU:N	1:D:452:GLU:OE1	2.28	0.67
1:E:454:ARG:C	1:E:455:ILE:HD12	2.19	0.67
1:C:644:HIS:HB3	1:C:653:TRP:HE1	1.60	0.67
1:A:542:HIS:HE1	1:A:568:ARG:HG2	1.59	0.67
1:A:450:LEU:O	1:A:738:PRO:CD	2.43	0.66
1:C:703:HIS:ND1	1:E:577:GLU:OE1	2.28	0.66
1:A:455:ILE:C	1:A:456:LEU:HG	2.19	0.66
1:D:450:LEU:O	1:D:738:PRO:HD2	1.95	0.66
1:C:692:ILE:CD1	1:C:706:LEU:HD12	2.25	0.66
1:D:712:PRO:CD	1:D:730:GLU:HG3	2.25	0.66
1:A:542:HIS:CE1	1:A:568:ARG:CG	2.79	0.65
1:D:452:GLU:N	1:D:452:GLU:CD	2.54	0.65
1:C:605:ARG:HG3	1:C:629:ASN:CG	2.20	0.65
1:A:607:GLY:O	1:A:626:PRO:CG	2.43	0.65
1:D:457:THR:CG2	1:D:486:TRP:CH2	2.79	0.65
1:D:538:VAL:C	1:D:539:PHE:CD1	2.74	0.65
1:B:628:GLU:HB3	1:B:646:SER:OG	1.96	0.65
1:C:731:ASP:OD1	1:C:733:THR:HG22	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:457:THR:HG23	1:D:458:THR:N	2.12	0.65
1:C:625:ALA:N	1:C:626:PRO:HD2	2.10	0.65
1:A:628:GLU:HB2	1:A:647:ASP:HB2	1.79	0.65
1:D:455:ILE:C	1:D:456:LEU:HG	2.20	0.65
1:B:615:ALA:CB	1:C:719:HIS:HE2	2.09	0.64
1:D:712:PRO:HD2	1:D:730:GLU:HG3	1.77	0.64
1:C:480:ASP:HB2	1:C:482:LEU:CD1	2.26	0.64
1:D:650:VAL:HG13	1:D:664:PHE:HD2	1.61	0.64
1:A:457:THR:HG1	1:A:486:TRP:HH2	1.44	0.64
1:D:456:LEU:O	1:D:457:THR:HB	1.97	0.64
1:C:703:HIS:O	1:C:704:THR:OG1	2.11	0.64
1:D:650:VAL:CG1	1:D:664:PHE:HD2	2.10	0.64
1:C:536:ARG:HG2	1:C:537:ALA:H	1.62	0.64
1:B:494:LEU:N	1:B:494:LEU:HD12	2.12	0.63
1:A:589:TYR:CE2	1:A:605:ARG:HG3	2.33	0.63
1:C:480:ASP:HB2	1:C:482:LEU:HD11	1.79	0.63
1:C:689:ASP:OD1	1:C:691:THR:HG22	1.98	0.63
1:D:537:ALA:O	1:D:539:PHE:HE1	1.80	0.63
1:A:709:HIS:O	1:A:710:THR:OG1	2.12	0.63
1:C:563:ARG:HG2	1:C:587:TYR:HD2	1.63	0.63
1:D:450:LEU:HD12	1:D:450:LEU:H	1.64	0.63
1:A:648:SER:C	1:A:649:THR:HG23	2.24	0.62
1:D:451:ASN:HD22	1:D:451:ASN:C	2.06	0.62
1:A:457:THR:OG1	1:A:486:TRP:HH2	1.82	0.62
1:C:451:ASN:O	1:C:738:PRO:HD3	2.00	0.62
1:C:470:GLY:HA3	2:C:838:HOH:O	1.99	0.62
1:C:602:SER:HB2	1:C:612:TRP:HZ3	1.65	0.62
1:C:643:VAL:CG1	1:C:683:LEU:HD21	2.29	0.62
1:D:652:LEU:HD12	1:D:700:GLN:NE2	2.15	0.62
1:D:735:ARG:NH1	1:D:735:ARG:HB2	2.15	0.62
1:C:692:ILE:HG12	1:C:706:LEU:HB2	1.80	0.61
1:C:664:PHE:CD1	1:C:700:GLN:HG3	2.35	0.61
1:C:692:ILE:CD1	1:C:706:LEU:HB2	2.31	0.61
1:C:692:ILE:CG2	1:C:713:VAL:HG11	2.29	0.61
1:C:602:SER:HB2	1:C:612:TRP:CZ3	2.35	0.61
1:A:601:ALA:CB	1:A:642:LEU:HD11	2.31	0.61
1:D:526:ARG:NH1	1:D:539:PHE:HE2	1.99	0.61
1:A:560:SER:HG	1:A:570:TRP:HZ3	1.48	0.60
1:C:692:ILE:HD11	1:C:706:LEU:HB2	1.82	0.60
1:E:647:ASP:HB2	1:E:670:TRP:CZ3	2.36	0.60
1:A:560:SER:CB	1:A:570:TRP:HZ3	2.14	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:457:THR:HG21	1:D:486:TRP:CH2	2.36	0.60
1:D:539:PHE:C	1:D:540:GLU:HG2	2.26	0.60
1:D:619:GLU:OE2	1:D:622:VAL:CG2	2.50	0.59
1:C:568:ARG:HB3	1:C:578:HIS:CD2	2.36	0.59
1:C:694:LEU:CD2	1:C:725:LEU:HD21	2.21	0.59
1:D:542:HIS:NE2	1:D:570:TRP:HH2	1.99	0.59
1:D:601:ALA:HB2	1:D:642:LEU:HD11	1.84	0.59
1:B:711:GLU:HG3	1:B:731:ASP:HB3	1.83	0.59
1:D:460:ARG:CZ	1:D:503:TRP:CZ2	2.86	0.59
1:A:473:LEU:HD11	1:A:494:LEU:CD2	2.32	0.59
1:C:480:ASP:HB2	1:C:482:LEU:HG	1.85	0.59
1:C:587:TYR:CD1	1:C:605:ARG:HD3	2.38	0.59
1:C:649:THR:HG21	1:C:651:HIS:CE1	2.37	0.59
1:D:455:ILE:HB	1:D:734:ILE:HB	1.84	0.59
1:D:539:PHE:N	1:D:539:PHE:HD1	2.01	0.59
1:A:538:VAL:CG1	1:A:539:PHE:N	2.63	0.58
1:A:456:LEU:C	1:A:457:THR:HG23	2.28	0.58
1:D:457:THR:HG22	1:D:486:TRP:CH2	2.37	0.58
1:A:542:HIS:HE1	1:A:568:ARG:CG	2.15	0.58
1:C:616:THR:CG2	1:C:618:LYS:HB2	2.34	0.57
1:A:537:ALA:O	1:A:539:PHE:CE1	2.57	0.57
1:B:463:VAL:HG12	1:B:714:HIS:HD2	1.69	0.57
1:A:730:GLU:HB3	2:A:837:HOH:O	2.04	0.57
1:C:589:TYR:OH	1:C:629:ASN:ND2	2.37	0.57
1:D:574:THR:OG1	1:D:576:THR:HG23	2.04	0.57
1:D:460:ARG:CG	1:D:461:GLU:H	2.15	0.57
1:A:456:LEU:O	1:A:457:THR:CB	2.52	0.57
1:C:481:LYS:HD3	1:C:481:LYS:N	2.20	0.57
1:C:670:TRP:CZ3	1:D:545:TYR:OH	2.45	0.57
1:D:712:PRO:CG	1:D:730:GLU:HG3	2.35	0.57
1:D:661:LEU:O	1:D:662:HIS:HB2	2.04	0.57
1:E:689:ASP:O	1:E:690:ARG:HB2	2.04	0.57
1:A:537:ALA:O	1:A:539:PHE:HE1	1.87	0.56
1:A:721:GLU:C	1:A:721:GLU:OE1	2.48	0.56
1:C:590:ALA:HB1	1:C:633:LEU:HD22	1.85	0.56
1:C:599:MET:CE	1:C:620:ARG:HH11	2.18	0.56
1:C:664:PHE:HD1	1:C:700:GLN:HG3	1.70	0.56
1:E:647:ASP:O	1:E:648:SER:HB2	2.05	0.56
1:A:560:SER:OG	1:A:570:TRP:HZ3	1.88	0.56
1:A:453:PRO:HD2	1:A:736:ILE:CG2	2.35	0.56
1:C:584:HIS:O	1:C:585:THR:HG22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:631:VAL:HG11	1:E:672:ARG:HG2	1.86	0.56
1:D:558:VAL:HG23	1:D:572:VAL:HG12	1.87	0.56
1:B:578:HIS:CE1	1:C:472:SER:HB2	2.41	0.56
1:C:586:ASP:HB3	1:C:605:ARG:HB3	1.87	0.56
1:C:616:THR:HG23	1:C:618:LYS:CB	2.35	0.56
1:A:536:ARG:CA	1:A:536:ARG:HE	2.18	0.56
1:A:542:HIS:HD1	1:A:568:ARG:CZ	2.19	0.56
1:A:625:ALA:N	1:A:626:PRO:CD	2.69	0.56
1:D:643:VAL:HG11	1:D:683:LEU:CD2	2.36	0.56
1:C:480:ASP:HB2	1:C:482:LEU:CG	2.35	0.56
1:C:481:LYS:HD2	1:C:502:ASP:HA	1.88	0.56
1:C:536:ARG:CG	1:C:537:ALA:H	2.18	0.56
1:C:643:VAL:HG22	1:C:652:LEU:HD12	1.88	0.56
1:A:589:TYR:HE2	1:A:605:ARG:HG3	1.70	0.56
1:C:460:ARG:HD3	1:C:480:ASP:CG	2.31	0.56
1:A:715:SER:OG	1:A:728:ALA:HB3	2.06	0.55
1:C:632:SER:HB3	1:C:674:VAL:HG22	1.87	0.55
1:C:643:VAL:CG2	1:C:683:LEU:HD21	2.37	0.55
1:C:692:ILE:HG13	1:C:706:LEU:HD12	1.88	0.55
1:A:568:ARG:HG2	1:A:568:ARG:NH2	2.22	0.55
1:C:563:ARG:HG2	1:C:587:TYR:CD2	2.40	0.55
1:C:584:HIS:HB3	1:C:606:ASP:OD2	2.06	0.55
1:C:724:THR:HB	1:C:737:TRP:O	2.07	0.55
1:D:453:PRO:CD	1:D:453:PRO:O	2.55	0.55
1:D:479:GLY:HA2	1:D:503:TRP:CD2	2.42	0.55
1:D:547:LEU:HD11	1:D:587:TYR:CD2	2.41	0.55
1:A:536:ARG:CA	1:A:536:ARG:NE	2.68	0.55
1:A:539:PHE:N	1:A:539:PHE:CD1	2.73	0.55
1:A:724:THR:CG2	1:A:736:ILE:HG12	2.37	0.55
1:C:628:GLU:HG2	1:C:629:ASN:N	2.08	0.55
1:B:631:VAL:HG11	1:B:672:ARG:HG2	1.89	0.54
1:D:452:GLU:HB2	1:D:453:PRO:HD3	1.89	0.54
1:D:558:VAL:CG2	1:D:572:VAL:HG12	2.36	0.54
1:D:657:SER:HB3	1:D:659:GLU:HG2	1.89	0.54
1:D:712:PRO:HG2	1:D:730:GLU:HG3	1.89	0.54
1:C:481:LYS:HD2	1:C:502:ASP:CA	2.37	0.54
1:C:578:HIS:ND1	1:C:578:HIS:N	2.56	0.54
1:E:480:ASP:OD2	1:E:482:LEU:HD12	2.06	0.54
1:A:463:VAL:HG12	1:A:714:HIS:HD2	1.72	0.54
1:A:560:SER:HB2	1:A:570:TRP:HZ3	1.70	0.54
1:A:650:VAL:CG2	1:A:671:VAL:HG11	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:650:VAL:HB	1:C:664:PHE:HB2	1.88	0.54
1:D:456:LEU:C	1:D:457:THR:HG22	2.31	0.54
1:D:650:VAL:CG1	1:D:664:PHE:CD2	2.88	0.54
1:A:643:VAL:CG2	1:A:676:PHE:CE2	2.90	0.54
1:A:650:VAL:HG22	1:A:671:VAL:HG11	1.90	0.54
1:C:460:ARG:O	1:C:732:GLY:HA2	2.07	0.54
1:C:599:MET:CE	1:C:620:ARG:HD2	2.19	0.54
1:C:601:ALA:HA	1:C:610:ARG:O	2.08	0.54
1:A:456:LEU:O	1:A:457:THR:OG1	2.26	0.54
1:E:644:HIS:CD2	1:E:653:TRP:HE1	2.23	0.54
1:D:625:ALA:N	1:D:626:PRO:CD	2.72	0.53
1:D:628:GLU:HB2	1:D:647:ASP:HB2	1.91	0.53
1:A:465:VAL:HG13	1:A:728:ALA:HB2	1.90	0.53
1:A:589:TYR:HE2	1:A:605:ARG:CG	2.21	0.53
1:D:520:SER:O	1:D:546:VAL:HG23	2.07	0.53
1:A:526:ARG:CZ	1:A:539:PHE:HE2	2.20	0.53
1:A:536:ARG:NE	1:A:536:ARG:HA	2.23	0.53
1:A:711:GLU:HB2	1:A:730:GLU:OE1	2.08	0.53
1:A:729:SER:OG	1:A:733:THR:OG1	2.26	0.53
1:D:661:LEU:HD23	1:D:700:GLN:HE21	1.73	0.53
1:C:569:LEU:N	1:C:569:LEU:HD12	2.23	0.53
1:C:460:ARG:CD	1:C:478:SER:OG	2.55	0.53
1:D:707:GLU:HB3	1:D:709:HIS:CD2	2.44	0.53
1:C:563:ARG:HA	1:C:587:TYR:HB3	1.91	0.53
1:E:643:VAL:HG11	1:E:683:LEU:HD11	1.91	0.53
1:C:591:VAL:C	1:C:633:LEU:HD21	2.34	0.53
1:D:643:VAL:HG11	1:D:683:LEU:CD1	2.38	0.53
1:D:717:ALA:O	1:D:725:LEU:HD12	2.09	0.53
1:E:643:VAL:HG21	1:E:683:LEU:HD21	1.91	0.53
1:A:560:SER:HB2	1:A:570:TRP:CZ3	2.44	0.52
1:D:455:ILE:C	1:D:456:LEU:CG	2.80	0.52
1:D:568:ARG:HB2	1:D:568:ARG:NH2	2.24	0.52
1:C:643:VAL:HG11	1:C:683:LEU:HD21	1.91	0.52
1:D:542:HIS:CD2	1:D:546:VAL:HG13	2.45	0.52
1:A:724:THR:HG22	1:A:725:LEU:N	2.23	0.52
1:A:643:VAL:CG2	1:A:676:PHE:CZ	2.93	0.52
1:C:557:MET:HE1	1:C:614:VAL:HG11	1.92	0.52
1:E:609:ILE:HB	1:E:623:LEU:HB2	1.92	0.52
1:A:542:HIS:CE1	1:A:568:ARG:NH2	2.78	0.52
1:D:560:SER:HB2	1:D:570:TRP:HZ3	1.73	0.52
1:C:692:ILE:HG22	1:C:713:VAL:HG11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:716:VAL:HG13	1:E:725:LEU:HD11	1.92	0.52
1:A:711:GLU:CB	1:A:730:GLU:OE1	2.58	0.52
1:C:599:MET:HE2	1:C:620:ARG:HD3	1.89	0.52
1:C:651:HIS:CE1	2:C:802:HOH:O	2.62	0.52
1:D:461:GLU:OE2	1:D:730:GLU:CG	2.58	0.52
1:D:657:SER:CB	1:D:659:GLU:HG2	2.40	0.52
1:A:601:ALA:HB2	1:A:642:LEU:HD11	1.91	0.52
1:C:589:TYR:HE2	1:C:629:ASN:HB3	1.75	0.51
1:D:450:LEU:O	1:D:737:TRP:HB3	2.10	0.51
1:C:536:ARG:HG2	1:C:537:ALA:N	2.25	0.51
1:D:623:LEU:HB3	1:D:653:TRP:CE3	2.45	0.51
1:E:692:ILE:HD12	1:E:727:SER:HB3	1.92	0.51
1:C:482:LEU:HD13	1:C:484:HIS:CD2	2.45	0.51
1:C:567:ALA:HB3	1:C:581:LEU:HB2	1.92	0.51
1:D:461:GLU:CG	1:D:503:TRP:HH2	2.19	0.51
1:A:644:HIS:CE1	1:A:646:SER:HB2	2.46	0.51
1:C:562:SER:N	1:C:588:VAL:HG23	2.25	0.51
1:D:456:LEU:O	1:D:457:THR:CB	2.59	0.51
1:C:692:ILE:HG13	1:C:706:LEU:HB2	1.89	0.51
1:D:724:THR:HG23	1:D:737:TRP:O	2.11	0.51
1:D:735:ARG:HB2	1:D:735:ARG:HH11	1.75	0.51
1:A:609:ILE:HD12	1:A:630:VAL:HG21	1.92	0.50
1:A:579:ALA:HB2	2:A:834:HOH:O	2.10	0.50
1:C:652:LEU:HD21	1:C:697:VAL:HG13	1.93	0.50
1:A:723:THR:O	1:A:738:PRO:HA	2.11	0.50
1:C:643:VAL:CG2	1:C:683:LEU:HD11	2.25	0.50
1:E:557:MET:HE1	1:E:578:HIS:CG	2.45	0.50
1:C:605:ARG:CG	1:C:629:ASN:ND2	2.54	0.50
1:D:506:ALA:CB	1:D:548:ASP:HA	2.42	0.50
1:D:624:GLN:O	1:D:625:ALA:HB2	2.12	0.50
1:E:610:ARG:HG2	1:E:622:VAL:HG22	1.92	0.50
1:A:473:LEU:HD11	1:A:494:LEU:HD22	1.93	0.50
1:D:652:LEU:HB2	1:D:700:GLN:HE22	1.76	0.50
1:D:548:ASP:OD2	1:D:591:VAL:HG22	2.11	0.50
1:D:643:VAL:HG11	1:D:683:LEU:HD22	1.93	0.50
1:A:737:TRP:N	1:A:737:TRP:CD1	2.80	0.50
1:C:613:ASP:OD2	1:C:616:THR:HG22	2.12	0.50
1:E:527:LEU:HD13	1:E:539:PHE:HE2	1.76	0.50
1:C:643:VAL:HG21	1:C:683:LEU:HD21	1.93	0.50
1:C:527:LEU:O	1:C:536:ARG:HB3	2.11	0.49
1:D:542:HIS:CE1	1:D:568:ARG:NH2	2.74	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:625:ALA:H	1:A:626:PRO:HD2	1.77	0.49
1:C:643:VAL:HG21	1:C:683:LEU:CD2	2.42	0.49
1:C:682:LEU:HD23	1:C:695:TRP:O	2.12	0.49
1:A:568:ARG:HB2	1:A:570:TRP:CZ3	2.48	0.49
1:D:537:ALA:O	1:D:539:PHE:CE1	2.63	0.49
1:D:600:VAL:HG23	1:D:614:VAL:HG12	1.94	0.49
1:E:656:ALA:O	1:E:657:SER:CB	2.60	0.49
1:A:545:TYR:H	1:A:545:TYR:HD1	1.54	0.49
1:D:453:PRO:HD2	1:D:453:PRO:O	2.13	0.49
1:C:458:THR:OG1	1:C:460:ARG:NH1	2.46	0.49
1:C:664:PHE:HD1	1:C:700:GLN:CG	2.21	0.49
1:A:662:HIS:ND1	1:A:700:GLN:HB3	2.28	0.49
1:D:526:ARG:CZ	1:D:539:PHE:HE2	2.25	0.49
1:D:643:VAL:HG21	1:D:683:LEU:HD11	1.95	0.49
1:A:545:TYR:CD1	1:A:545:TYR:N	2.72	0.49
1:A:609:ILE:HG22	1:A:623:LEU:HD12	1.94	0.49
1:C:625:ALA:N	1:C:626:PRO:CD	2.75	0.49
1:A:526:ARG:CZ	1:A:539:PHE:CE2	2.95	0.48
1:B:563:ARG:NH2	1:B:587:TYR:OH	2.46	0.48
1:C:692:ILE:HG13	1:C:706:LEU:CG	2.42	0.48
1:A:455:ILE:C	1:A:456:LEU:CG	2.82	0.48
1:B:644:HIS:CE1	1:B:651:HIS:HB2	2.48	0.48
1:D:545:TYR:H	1:D:545:TYR:HD1	1.56	0.48
1:B:453:PRO:CB	1:B:736:ILE:O	2.61	0.48
1:D:662:HIS:HB3	1:D:700:GLN:NE2	2.24	0.48
1:C:673:ALA:CB	2:C:834:HOH:O	2.56	0.48
1:D:545:TYR:CD1	1:D:545:TYR:N	2.78	0.48
1:A:453:PRO:CB	2:A:857:HOH:O	2.62	0.48
1:C:601:ALA:O	1:C:633:LEU:HD11	2.14	0.48
1:A:548:ASP:HB3	1:A:591:VAL:HG22	1.95	0.47
1:A:644:HIS:O	1:A:650:VAL:HG13	2.14	0.47
1:A:724:THR:HG21	1:A:736:ILE:HG12	1.96	0.47
1:E:563:ARG:NH2	1:E:587:TYR:OH	2.46	0.47
1:C:525:VAL:HB	1:C:539:PHE:HB2	1.95	0.47
1:A:677:SER:OG	1:A:679:ASP:OD1	2.28	0.47
1:C:649:THR:CG2	1:C:651:HIS:CE1	2.98	0.47
1:D:453:PRO:HG3	1:D:736:ILE:HD11	1.97	0.47
1:D:623:LEU:HD13	1:D:653:TRP:CB	2.44	0.47
1:D:643:VAL:HG11	1:D:683:LEU:HD13	1.97	0.47
1:A:479:GLY:HA3	1:A:503:TRP:CE3	2.50	0.47
1:A:487:ASP:CG	1:A:489:ALA:O	2.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:485:VAL:HG13	1:C:494:LEU:HB2	1.95	0.47
1:C:674:VAL:HA	1:C:684:ALA:O	2.14	0.47
1:D:453:PRO:O	1:D:454:ARG:HD2	2.06	0.47
1:E:692:ILE:HD13	1:E:713:VAL:HG11	1.96	0.47
1:A:463:VAL:HG11	1:A:505:ARG:NH1	2.30	0.47
1:C:692:ILE:CG1	1:C:706:LEU:HD12	2.44	0.47
1:D:460:ARG:NH2	1:D:503:TRP:HZ2	2.13	0.47
1:E:626:PRO:O	1:E:627:ALA:HB3	2.15	0.47
1:E:692:ILE:HB	1:E:706:LEU:HB2	1.96	0.47
1:D:451:ASN:HA	1:D:737:TRP:HA	1.97	0.47
1:E:719:HIS:CE1	1:E:721:GLU:HG2	2.50	0.47
1:A:479:GLY:CA	1:A:503:TRP:CE3	2.98	0.46
1:A:644:HIS:HB3	1:A:653:TRP:HE1	1.80	0.46
1:A:730:GLU:HG2	1:A:731:ASP:HB2	1.97	0.46
1:D:623:LEU:HD13	1:D:653:TRP:HB3	1.97	0.46
1:D:704:THR:HG22	1:D:705:THR:N	2.31	0.46
1:C:685:SER:OG	1:C:695:TRP:NE1	2.44	0.46
1:A:522:ASP:O	1:A:523:ALA:HB3	2.15	0.46
1:D:724:THR:HG22	1:D:725:LEU:N	2.31	0.46
1:E:547:LEU:HD11	1:E:563:ARG:HG3	1.98	0.46
1:A:473:LEU:HD11	1:A:494:LEU:HD21	1.97	0.46
1:C:536:ARG:CG	1:C:537:ALA:N	2.79	0.46
1:C:562:SER:O	1:C:588:VAL:HG22	2.16	0.46
1:E:616:THR:OG1	1:E:618:LYS:HB2	2.16	0.46
1:B:587:TYR:N	1:B:587:TYR:CD1	2.84	0.46
1:C:644:HIS:ND1	1:C:651:HIS:HB2	2.31	0.46
1:D:726:ALA:HA	1:D:735:ARG:O	2.15	0.46
1:D:542:HIS:CD2	1:D:546:VAL:CG1	2.99	0.46
1:C:625:ALA:HB3	1:C:653:TRP:HH2	1.80	0.46
1:D:563:ARG:CB	1:D:587:TYR:CZ	2.98	0.46
1:E:562:SER:OG	1:E:564:ASP:OD1	2.34	0.46
1:C:692:ILE:HD12	1:C:706:LEU:HD12	1.97	0.45
1:D:542:HIS:ND1	1:D:568:ARG:NH2	2.63	0.45
1:D:650:VAL:HG13	1:D:664:PHE:HB2	1.97	0.45
1:E:536:ARG:O	1:E:536:ARG:HG2	2.15	0.45
1:A:624:GLN:O	1:A:625:ALA:HB2	2.16	0.45
1:A:657:SER:OG	1:A:659:GLU:HG2	2.16	0.45
1:C:539:PHE:HE2	1:C:558:VAL:HG11	1.81	0.45
1:D:453:PRO:C	1:D:454:ARG:CD	2.62	0.45
1:A:458:THR:HG22	1:A:732:GLY:HA3	1.98	0.45
1:A:721:GLU:OE1	1:A:723:THR:HB	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:569:LEU:N	1:C:569:LEU:CD1	2.79	0.45
1:C:583:GLY:O	1:C:585:THR:N	2.49	0.45
1:D:737:TRP:CD1	1:D:737:TRP:N	2.85	0.45
1:A:457:THR:OG1	1:A:486:TRP:CH2	2.67	0.45
1:C:568:ARG:HB3	1:C:578:HIS:HD2	1.78	0.45
1:C:609:ILE:HB	1:C:623:LEU:HB2	1.98	0.45
1:D:719:HIS:CE1	1:D:721:GLU:HG2	2.51	0.45
1:E:527:LEU:HB3	1:E:536:ARG:HD3	1.98	0.45
1:E:628:GLU:HB3	1:E:646:SER:OG	2.16	0.45
1:C:703:HIS:C	1:C:704:THR:HG1	2.17	0.45
1:B:711:GLU:HB3	1:B:712:PRO:HD2	1.97	0.45
1:D:616:THR:O	1:D:618:LYS:N	2.47	0.45
1:D:696:ASP:O	1:D:700:GLN:N	2.49	0.45
1:C:458:THR:O	1:C:459:ASP:HB2	2.17	0.45
1:C:605:ARG:CG	1:C:629:ASN:CG	2.88	0.45
1:D:461:GLU:OE2	1:D:730:GLU:CB	2.64	0.45
1:D:644:HIS:CB	1:D:653:TRP:HE1	2.27	0.45
1:A:456:LEU:C	1:A:457:THR:OG1	2.60	0.45
1:A:620:ARG:HD2	1:A:620:ARG:HA	1.81	0.45
1:B:711:GLU:HB3	1:B:712:PRO:CD	2.47	0.45
1:C:688:ASP:HA	1:C:712:PRO:HB3	1.98	0.45
1:D:451:ASN:HB3	1:D:737:TRP:CE3	2.52	0.45
1:D:570:TRP:NE1	1:D:577:GLU:HB2	2.32	0.45
1:A:609:ILE:HG13	1:A:633:LEU:HD11	1.99	0.45
1:D:712:PRO:HG2	1:D:730:GLU:CG	2.46	0.45
1:A:719:HIS:HB2	1:A:724:THR:HB	1.99	0.45
1:A:724:THR:CG2	1:A:725:LEU:N	2.79	0.45
1:B:487:ASP:N	1:B:494:LEU:CD1	2.80	0.45
1:C:649:THR:HG21	1:C:651:HIS:HE1	1.81	0.45
1:D:633:LEU:HD12	1:D:633:LEU:C	2.42	0.45
1:B:483:ILE:HB	1:B:497:LEU:HB2	1.99	0.44
1:C:637:PRO:HG3	1:C:678:PRO:O	2.16	0.44
1:C:453:PRO:N	1:C:453:PRO:C	2.64	0.44
1:D:460:ARG:NE	1:D:503:TRP:HZ2	2.16	0.44
1:A:685:SER:O	1:A:692:ILE:HA	2.17	0.44
1:C:547:LEU:HD12	1:C:589:TYR:HD1	1.82	0.44
1:E:455:ILE:N	1:E:455:ILE:CD1	2.70	0.44
1:A:641:MET:HE3	1:A:697:VAL:CG1	2.47	0.44
1:B:685:SER:O	1:B:692:ILE:HA	2.18	0.44
1:D:502:ASP:HB3	1:D:521:ASP:HB2	2.00	0.44
1:D:652:LEU:HB2	1:D:700:GLN:NE2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:711:GLU:HG2	1:E:731:ASP:HB3	2.00	0.44
1:A:551:PHE:CE1	1:A:558:VAL:CG2	3.01	0.44
1:A:568:ARG:CG	1:A:568:ARG:NH2	2.80	0.44
1:A:662:HIS:HB2	1:A:700:GLN:HG2	2.00	0.44
1:D:699:ALA:O	1:D:700:GLN:CB	2.60	0.43
1:E:525:VAL:HB	1:E:539:PHE:HB2	1.99	0.43
1:A:482:LEU:HD22	1:A:496:THR:CG2	2.48	0.43
1:A:536:ARG:HE	1:A:536:ARG:HA	1.80	0.43
1:E:631:VAL:O	1:E:631:VAL:HG12	2.16	0.43
1:E:662:HIS:CG	1:E:663:THR:N	2.86	0.43
1:A:450:LEU:O	1:A:738:PRO:HD2	2.17	0.43
1:D:609:ILE:HD11	1:D:630:VAL:HG11	2.00	0.43
1:A:539:PHE:N	1:A:539:PHE:HD1	2.15	0.43
1:A:719:HIS:CG	1:A:720:PRO:HD2	2.53	0.43
1:C:460:ARG:HD3	1:C:480:ASP:OD2	2.18	0.43
1:C:501:THR:HG22	1:C:501:THR:O	2.18	0.43
1:C:605:ARG:CG	1:C:629:ASN:HD21	2.12	0.43
1:D:571:ASN:OD1	1:D:574:THR:HG22	2.18	0.43
1:A:674:VAL:HA	1:A:684:ALA:O	2.18	0.43
1:D:619:GLU:OE2	1:D:622:VAL:HG22	2.16	0.43
1:D:538:VAL:CG1	1:D:539:PHE:H	2.30	0.43
1:D:631:VAL:HG12	1:D:631:VAL:O	2.19	0.43
1:A:558:VAL:HG23	1:A:572:VAL:HG22	2.00	0.43
1:D:454:ARG:HG3	1:D:735:ARG:CG	2.47	0.43
1:E:726:ALA:HA	1:E:735:ARG:O	2.19	0.43
1:B:481:LYS:HG2	1:B:502:ASP:C	2.44	0.43
1:D:589:TYR:CE1	1:D:605:ARG:HG3	2.54	0.43
1:D:600:VAL:HG23	1:D:614:VAL:CG1	2.48	0.43
1:A:482:LEU:HD22	1:A:496:THR:HG21	2.01	0.42
1:A:625:ALA:HB3	1:A:653:TRP:HH2	1.84	0.42
1:D:457:THR:CG2	1:D:458:THR:N	2.79	0.42
1:A:538:VAL:C	1:A:539:PHE:CD1	2.97	0.42
1:A:542:HIS:HD1	1:A:568:ARG:NH2	2.16	0.42
1:E:664:PHE:HB3	1:E:695:TRP:CZ3	2.54	0.42
1:C:460:ARG:HD3	1:C:480:ASP:OD1	2.18	0.42
1:D:471:GLY:HA3	1:D:472:SER:CB	2.49	0.42
1:D:598:SER:O	1:D:614:VAL:HG22	2.19	0.42
1:A:479:GLY:HA2	1:A:503:TRP:CD2	2.55	0.42
1:A:548:ASP:CG	1:A:590:ALA:HA	2.45	0.42
1:B:647:ASP:O	1:B:648:SER:HB2	2.20	0.42
1:C:482:LEU:HD12	1:C:482:LEU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:664:PHE:CE1	1:C:700:GLN:CG	3.02	0.42
1:D:647:ASP:CB	1:D:670:TRP:HD1	2.33	0.42
1:D:487:ASP:OD2	1:D:490:SER:HB3	2.19	0.42
1:E:454:ARG:CA	1:E:455:ILE:HD12	2.49	0.42
1:E:647:ASP:HB2	1:E:670:TRP:CH2	2.55	0.42
1:A:542:HIS:ND1	1:A:568:ARG:NH2	2.68	0.42
1:C:701:GLU:HG3	1:E:579:ALA:HB1	2.00	0.42
1:D:462:ALA:O	1:D:728:ALA:HB1	2.19	0.42
1:E:562:SER:OG	1:E:564:ASP:CG	2.62	0.42
1:A:589:TYR:CD2	1:A:605:ARG:HG3	2.54	0.42
1:A:647:ASP:OD1	1:A:670:TRP:CD2	2.73	0.42
1:D:461:GLU:OE2	1:D:730:GLU:HA	2.19	0.42
1:A:726:ALA:HA	1:A:735:ARG:O	2.20	0.42
1:B:494:LEU:N	1:B:494:LEU:CD1	2.81	0.42
1:A:452:GLU:HB3	1:A:736:ILE:CG2	2.49	0.42
1:B:494:LEU:HD12	1:B:494:LEU:H	1.81	0.42
1:B:566:THR:HG21	1:B:568:ARG:HH21	1.84	0.42
1:C:480:ASP:CB	1:C:482:LEU:HG	2.48	0.42
1:C:624:GLN:O	1:C:625:ALA:HB2	2.20	0.42
1:E:453:PRO:CB	1:E:736:ILE:O	2.68	0.42
1:E:522:ASP:O	1:E:523:ALA:HB3	2.19	0.42
1:A:502:ASP:HB3	1:A:521:ASP:HB2	2.02	0.41
1:B:527:LEU:HD12	1:B:527:LEU:HA	1.92	0.41
1:C:692:ILE:HD11	1:C:706:LEU:CB	2.48	0.41
1:D:560:SER:CB	1:D:570:TRP:HZ3	2.33	0.41
1:D:602:SER:CA	1:D:633:LEU:HD21	2.50	0.41
1:D:609:ILE:HB	1:D:623:LEU:HB2	2.02	0.41
1:A:453:PRO:HG2	1:A:736:ILE:HG21	2.02	0.41
1:A:525:VAL:HG23	1:A:546:VAL:HG21	2.02	0.41
1:C:719:HIS:ND1	1:C:721:GLU:HG2	2.34	0.41
1:C:733:THR:HG21	1:C:735:ARG:HH21	1.86	0.41
1:A:453:PRO:HD2	1:A:736:ILE:HG21	2.03	0.41
1:A:609:ILE:CD1	1:A:630:VAL:HG21	2.50	0.41
1:D:644:HIS:O	1:D:650:VAL:HA	2.21	0.41
1:B:529:ASP:HB3	1:B:534:GLU:HG2	2.02	0.41
1:C:607:GLY:O	1:C:626:PRO:CG	2.64	0.41
1:E:606:ASP:OD1	1:E:608:THR:OG1	2.25	0.41
1:B:692:ILE:HB	1:B:706:LEU:HB2	2.02	0.41
1:C:450:LEU:CB	1:C:704:THR:HG23	2.50	0.41
1:C:692:ILE:HG13	1:C:706:LEU:CB	2.50	0.41
1:D:675:ALA:O	1:D:683:LEU:HD12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:736:ILE:HD13	1:E:736:ILE:HG21	1.89	0.41
1:B:522:ASP:OD1	1:B:524:THR:HB	2.20	0.41
1:B:682:LEU:HD23	1:B:695:TRP:O	2.20	0.41
1:C:539:PHE:HB3	1:C:570:TRP:CZ3	2.56	0.41
1:D:542:HIS:HE1	1:D:568:ARG:HH21	1.61	0.41
1:D:692:ILE:HB	1:D:706:LEU:HB2	2.02	0.41
1:E:662:HIS:CD2	1:E:663:THR:N	2.89	0.41
1:A:458:THR:HG22	1:A:732:GLY:CA	2.50	0.41
1:C:473:LEU:CD2	1:C:494:LEU:CD1	2.99	0.41
1:D:568:ARG:HH21	1:D:568:ARG:HB2	1.86	0.41
1:C:562:SER:O	1:C:588:VAL:CG2	2.69	0.41
1:D:589:TYR:CE1	1:D:605:ARG:HD3	2.56	0.41
1:A:646:SER:OG	1:A:647:ASP:N	2.53	0.41
1:C:711:GLU:HG3	1:C:712:PRO:HD2	2.02	0.41
1:D:547:LEU:HD11	1:D:587:TYR:HD2	1.83	0.41
1:D:554:ASP:OD2	1:D:556:SER:OG	2.39	0.41
1:D:644:HIS:CE1	1:D:646:SER:HB2	2.55	0.41
1:A:455:ILE:O	1:A:456:LEU:HD23	2.20	0.41
1:C:646:SER:OG	1:C:647:ASP:N	2.54	0.41
1:C:647:ASP:HB3	1:C:670:TRP:CD1	2.56	0.41
1:D:685:SER:O	1:D:692:ILE:HA	2.21	0.41
1:C:706:LEU:HB3	1:C:737:TRP:CZ3	2.56	0.40
1:C:473:LEU:HD22	1:C:494:LEU:HD12	2.03	0.40
1:E:460:ARG:HG3	1:E:480:ASP:HB3	2.02	0.40
1:A:452:GLU:HB3	1:A:736:ILE:HG23	2.04	0.40
1:A:721:GLU:CD	1:A:723:THR:HB	2.46	0.40
1:D:625:ALA:HB3	1:D:653:TRP:HH2	1.85	0.40
1:D:650:VAL:HG23	1:D:674:VAL:HG11	2.04	0.40
1:E:699:ALA:O	1:E:700:GLN:HB2	2.22	0.40
1:A:644:HIS:HE1	1:A:649:THR:O	2.05	0.40
1:C:643:VAL:HG12	1:C:674:VAL:HG21	2.03	0.40
1:D:572:VAL:HG23	1:D:573:ALA:N	2.36	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	282/302 (93%)	268 (95%)	10 (4%)	4 (1%)	9	19
1	B	286/302 (95%)	273 (96%)	12 (4%)	1 (0%)	36	58
1	C	288/302 (95%)	270 (94%)	15 (5%)	3 (1%)	12	28
1	D	280/302 (93%)	258 (92%)	15 (5%)	7 (2%)	4	8
1	E	286/302 (95%)	273 (96%)	12 (4%)	1 (0%)	36	58
All	All	1422/1510 (94%)	1342 (94%)	64 (4%)	16 (1%)	11	25

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	452	GLU
1	A	625	ALA
1	C	625	ALA
1	D	453	PRO
1	D	457	THR
1	D	625	ALA
1	A	457	THR
1	A	710	THR
1	B	583	GLY
1	D	472	SER
1	D	709	HIS
1	E	657	SER
1	D	617	GLY
1	C	450	LEU
1	C	648	SER
1	D	452	GLU

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	219/235 (93%)	198 (90%)	21 (10%)	8	17
1	B	223/235 (95%)	209 (94%)	14 (6%)	16	36
1	C	221/235 (94%)	198 (90%)	23 (10%)	7	14
1	D	219/235 (93%)	189 (86%)	30 (14%)	3	7
1	E	222/235 (94%)	210 (95%)	12 (5%)	20	42
All	All	1104/1175 (94%)	1004 (91%)	100 (9%)	9	19

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	458	THR
1	A	473	LEU
1	A	493	GLU
1	A	516	LEU
1	A	539	PHE
1	A	543	THR
1	A	545	TYR
1	A	547	LEU
1	A	549	ILE
1	A	569	LEU
1	A	628	GLU
1	A	633	LEU
1	A	635	PHE
1	A	642	LEU
1	A	644	HIS
1	A	683	LEU
1	A	703	HIS
1	A	721	GLU
1	A	734	ILE
1	A	736	ILE
1	A	737	TRP
1	B	465	VAL
1	B	481	LYS
1	B	495	HIS
1	B	524	THR
1	B	527	LEU
1	B	551	PHE
1	B	572	VAL
1	B	582	LYS
1	B	587	TYR
1	B	624	GLN

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Mol	Chain	Res	Type
1	B	628	GLU
1	B	638	ASP
1	B	710	THR
1	B	724	THR
1	C	452	GLU
1	C	457	THR
1	C	482	LEU
1	C	515	LEU
1	C	564	ASP
1	C	566	THR
1	C	587	TYR
1	C	602	SER
1	C	622	VAL
1	C	632	SER
1	C	633	LEU
1	C	642	LEU
1	C	643	VAL
1	C	644	HIS
1	C	649	THR
1	C	676	PHE
1	C	683	LEU
1	C	685	SER
1	C	692	ILE
1	C	711	GLU
1	C	725	LEU
1	C	736	ILE
1	C	737	TRP
1	D	450	LEU
1	D	451	ASN
1	D	453	PRO
1	D	458	THR
1	D	482	LEU
1	D	501	THR
1	D	516	LEU
1	D	539	PHE
1	D	540	GLU
1	D	543	THR
1	D	544	HIS
1	D	545	TYR
1	D	546	VAL
1	D	554	ASP
1	D	578	HIS

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Mol	Chain	Res	Type
1	D	582	LYS
1	D	611	LEU
1	D	616	THR
1	D	633	LEU
1	D	642	LEU
1	D	644	HIS
1	D	648	SER
1	D	659	GLU
1	D	703	HIS
1	D	710	THR
1	D	711	GLU
1	D	734	ILE
1	D	735	ARG
1	D	736	ILE
1	D	737	TRP
1	E	493	GLU
1	E	544	HIS
1	E	551	PHE
1	E	613	ASP
1	E	618	LYS
1	E	624	GLN
1	E	644	HIS
1	E	646	SER
1	E	685	SER
1	E	692	ILE
1	E	716	VAL
1	E	724	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	544	HIS
1	A	644	HIS
1	A	703	HIS
1	A	714	HIS
1	B	544	HIS
1	B	629	ASN
1	B	700	GLN
1	C	495	HIS
1	C	571	ASN
1	C	629	ASN
1	C	651	HIS

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Mol	Chain	Res	Type
1	D	544	HIS
1	D	578	HIS
1	D	651	HIS
1	D	700	GLN
1	D	709	HIS
1	E	644	HIS
1	E	662	HIS

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	286/302 (94%)	-1.50	0 100 100	19, 37, 68, 101	0
1	B	288/302 (95%)	-1.59	0 100 100	20, 36, 60, 79	0
1	C	290/302 (96%)	-1.47	0 100 100	19, 38, 69, 93	0
1	D	284/302 (94%)	-1.47	0 100 100	22, 40, 69, 92	0
1	E	288/302 (95%)	-1.62	0 100 100	23, 40, 61, 85	0
All	All	1436/1510 (95%)	-1.53	0 100 100	19, 38, 67, 101	0

There are no RSRZ outliers to report.

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.