



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:45 AM UTC

PDB ID : 1SOT / pdb_00001sot
Title : Crystal Structure of the DegS stress sensor
Authors : Wilken, C.; Kitzing, K.; Kurzbauer, R.; Ehrmann, M.; Clausen, T.
Deposited on : 2004-03-15
Resolution : 2.30 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

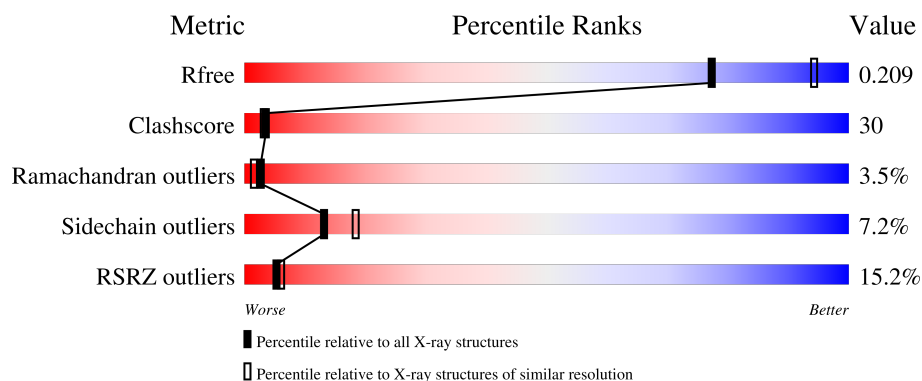
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	320	11% 49% 33% 6% 12%
1	B	320	15% 50% 36% 5% 8%
1	C	320	13% 53% 30% • • 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6617 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 Protease degS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	Se	0	0	0
			2066	1294	370	397	5			
1	B	294	Total	C	N	O	Se	91	0	0
			2175	1363	394	412	6			
1	C	281	Total	C	N	O	Se	91	0	0
			2046	1280	366	395	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	MSE	-	modified residue	UNP P31137
A	85	MSE	MET	modified residue	UNP P31137
A	213	MSE	MET	modified residue	UNP P31137
A	245	MSE	MET	modified residue	UNP P31137
A	319	MSE	MET	modified residue	UNP P31137
A	336	MSE	MET	modified residue	UNP P31137
A	356	HIS	-	expression tag	UNP P31137
A	357	HIS	-	expression tag	UNP P31137
A	358	HIS	-	expression tag	UNP P31137
A	359	HIS	-	expression tag	UNP P31137
A	360	HIS	-	expression tag	UNP P31137
A	361	HIS	-	expression tag	UNP P31137
B	42	MSE	-	modified residue	UNP P31137
B	85	MSE	MET	modified residue	UNP P31137
B	213	MSE	MET	modified residue	UNP P31137
B	245	MSE	MET	modified residue	UNP P31137
B	319	MSE	MET	modified residue	UNP P31137
B	336	MSE	MET	modified residue	UNP P31137
B	356	HIS	-	expression tag	UNP P31137
B	357	HIS	-	expression tag	UNP P31137
B	358	HIS	-	expression tag	UNP P31137
B	359	HIS	-	expression tag	UNP P31137
B	360	HIS	-	expression tag	UNP P31137

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Chain	Residue	Modelled	Actual	Comment	Reference
B	361	HIS	-	expression tag	UNP P31137
C	42	MSE	-	modified residue	UNP P31137
C	85	MSE	MET	modified residue	UNP P31137
C	213	MSE	MET	modified residue	UNP P31137
C	245	MSE	MET	modified residue	UNP P31137
C	319	MSE	MET	modified residue	UNP P31137
C	336	MSE	MET	modified residue	UNP P31137
C	356	HIS	-	expression tag	UNP P31137
C	357	HIS	-	expression tag	UNP P31137
C	358	HIS	-	expression tag	UNP P31137
C	359	HIS	-	expression tag	UNP P31137
C	360	HIS	-	expression tag	UNP P31137
C	361	HIS	-	expression tag	UNP P31137

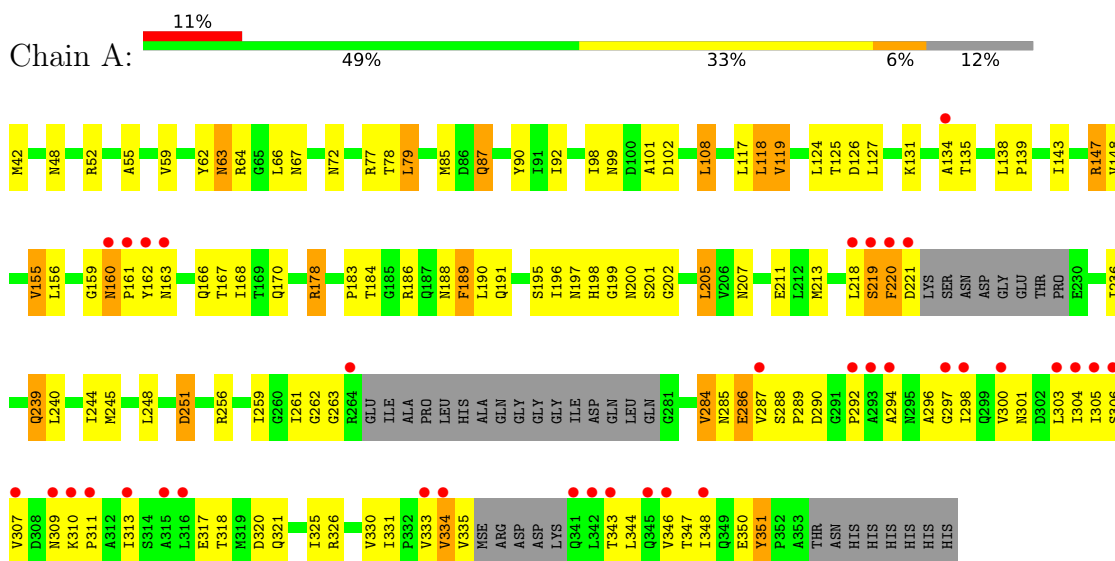
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	139	Total O 139 139	0	0
2	B	98	Total O 98 98	0	0
2	C	93	Total O 93 93	0	0

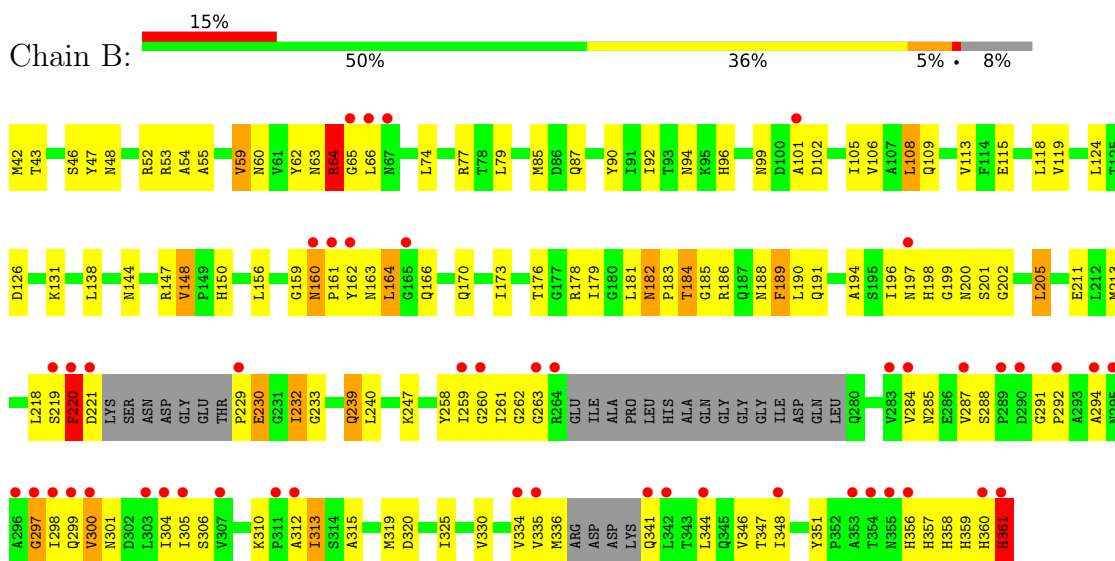
3 Residue-property plots [i](#)

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.

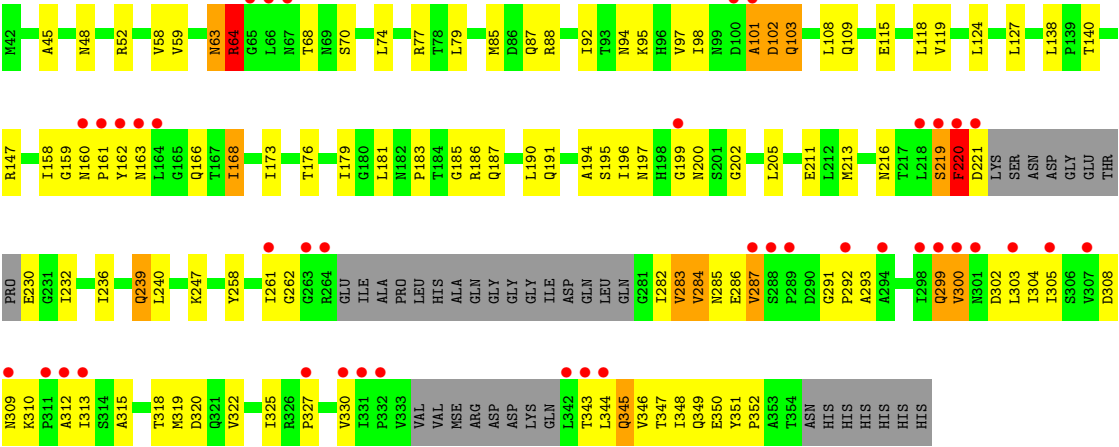
• Molecule 1: Protease degS



• Molecule 1: Protease degS



• Molecule 1: Protease degS



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	207.68Å 143.06Å 41.53Å 90.00° 90.07° 90.00°	Depositor
Resolution (Å)	15.00 – 2.30 15.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.30) 96.1 (15.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 1.95Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.198 , 0.248 0.208 , 0.209	Depositor DCC
R_{free} test set	4132 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 65.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6617	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% 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	0.89	1/2087 (0.0%)	1.09	9/2835 (0.3%)
1	B	0.98	3/2202 (0.1%)	1.19	18/2988 (0.6%)
1	C	0.81	3/2067 (0.1%)	1.96	22/2808 (0.8%)
All	All	0.89	7/6356 (0.1%)	1.46	49/8631 (0.6%)

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	B	0	1
1	C	0	1
All	All	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	77	ARG	C-N	-29.91	0.93	1.33
1	C	77	ARG	C-N	-17.54	1.11	1.33
1	B	64	ARG	C-N	-8.71	1.22	1.33
1	C	64	ARG	C-N	-6.68	1.23	1.33
1	B	361	HIS	C-OXT	6.47	1.36	1.23
1	C	310	LYS	N-CA	5.80	1.53	1.45
1	A	119	VAL	CA-CB	5.29	1.61	1.54

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	64	ARG	O-C-N	-83.59	28.57	123.03
1	B	77	ARG	CA-C-N	13.01	141.34	122.99
1	B	77	ARG	C-N-CA	13.01	141.34	122.99
1	B	194	ALA	N-CA-C	-9.71	95.90	110.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	77	ARG	CA-C-N	9.10	137.11	121.64
1	C	77	ARG	C-N-CA	9.10	137.11	121.64
1	B	64	ARG	CA-C-N	-9.08	112.84	121.46
1	B	64	ARG	C-N-CA	-9.08	112.84	121.46
1	B	77	ARG	O-C-N	-9.07	110.52	122.59
1	B	357	HIS	N-CA-C	8.07	120.08	111.28
1	C	77	ARG	O-C-N	-8.05	110.93	122.36
1	A	251	ASP	N-CA-C	-7.78	99.31	110.59
1	C	70	SER	CA-C-N	7.21	135.31	121.54
1	C	70	SER	C-N-CA	7.21	135.31	121.54
1	B	179	ILE	N-CA-C	6.90	118.73	112.17
1	C	68	THR	CA-C-N	-6.82	111.19	120.87
1	C	68	THR	C-N-CA	-6.82	111.19	120.87
1	C	70	SER	N-CA-C	6.70	125.06	110.80
1	C	115	GLU	N-CA-C	-6.69	99.68	110.32
1	A	189	PHE	N-CA-C	-6.56	101.07	110.59
1	B	115	GLU	N-CA-C	-6.39	100.16	110.32
1	B	189	PHE	N-CA-C	-6.08	101.77	110.59
1	A	87	GLN	N-CA-C	5.89	120.08	113.01
1	C	98	ILE	N-CA-C	5.89	119.57	112.80
1	B	43	THR	N-CA-C	-5.76	101.29	110.10
1	A	198	HIS	N-CA-C	5.73	118.30	111.71
1	B	54	ALA	N-CA-C	5.72	120.27	113.12
1	B	205	LEU	N-CA-C	-5.67	99.17	108.41
1	C	101	ALA	N-CA-C	5.57	117.81	111.02
1	C	45	ALA	N-CA-C	-5.51	102.85	110.35
1	B	176	THR	N-CA-C	5.50	118.20	111.82
1	B	182	ASN	CA-C-N	5.49	125.50	119.90
1	B	182	ASN	C-N-CA	5.49	125.50	119.90
1	C	68	THR	CB-CA-C	5.45	121.27	110.42
1	C	176	THR	N-CA-C	5.42	119.15	112.54
1	A	139	PRO	N-CA-C	-5.33	102.21	111.32
1	C	59	VAL	N-CA-C	5.31	116.50	108.96
1	C	68	THR	N-CA-C	-5.24	99.64	110.80
1	A	202	GLY	N-CA-C	-5.21	108.44	115.21
1	B	59	VAL	N-CA-C	5.16	116.61	109.29
1	C	68	THR	CA-C-O	-5.14	113.16	120.51
1	A	102	ASP	N-CA-C	-5.13	107.19	113.50
1	C	320	ASP	N-CA-C	-5.12	105.89	111.82
1	B	46	SER	N-CA-C	5.03	116.71	108.76
1	C	88	ARG	N-CA-C	-5.02	107.11	113.18
1	C	283	VAL	N-CA-C	5.01	114.49	107.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	LEU	CA-C-N	-5.01	114.44	121.55
1	A	118	LEU	C-N-CA	-5.01	114.44	121.55
1	C	70	SER	O-C-N	5.00	129.25	122.59

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	64	ARG	Mainchain
1	C	64	ARG	Mainchain

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	2066	0	2071	128	0
1	B	2175	0	2182	144	1
1	C	2046	0	2040	97	0
2	A	139	0	0	14	0
2	B	98	0	0	5	0
2	C	93	0	0	5	0
All	All	6617	0	6293	362	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 30.

All (362) 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:65:GLY:HA2	1:B:102:ASP:OD2	1.36	1.22
1:C:330:VAL:HG22	1:C:347:THR:HG22	1.23	1.11
1:C:185:GLY:HA2	1:C:221:ASP:HB2	1.17	1.10
1:B:184:THR:HG21	1:B:188:ASN:HB2	1.33	1.07
1:B:181:LEU:HD11	1:B:247:LYS:HE3	1.41	1.03
1:B:65:GLY:CA	1:B:102:ASP:OD2	2.08	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:PRO:HB3	1:A:199:GLY:HA3	1.47	0.97
1:A:330:VAL:HG22	1:A:347:THR:HG22	1.47	0.94
1:C:63:ASN:HD21	1:C:101:ALA:HA	1.32	0.93
1:B:305:ILE:HD12	1:B:334:VAL:HG12	1.51	0.93
1:B:356:HIS:ND1	1:B:361:HIS:CE1	2.38	0.92
1:A:239:GLN:H	1:A:239:GLN:HE21	1.07	0.91
1:B:184:THR:HG21	1:B:188:ASN:HD22	1.34	0.91
1:B:239:GLN:HE21	1:B:239:GLN:H	1.15	0.89
1:A:178:ARG:HH11	1:A:178:ARG:HB3	1.34	0.88
1:B:173:ILE:HD11	1:B:213:MSE:HE3	1.53	0.87
1:B:184:THR:CG2	1:B:188:ASN:HD22	1.86	0.87
1:C:63:ASN:ND2	1:C:101:ALA:HA	1.88	0.87
1:C:239:GLN:H	1:C:239:GLN:HE21	1.15	0.86
1:C:63:ASN:C	1:C:63:ASN:HD22	1.81	0.86
1:C:161:PRO:HB2	1:C:163:ASN:OD1	1.77	0.84
1:C:185:GLY:HA2	1:C:221:ASP:CB	2.04	0.84
1:A:263:GLY:HA2	1:A:285:ASN:HB2	1.61	0.82
1:B:356:HIS:CG	1:B:361:HIS:HE1	1.98	0.82
1:B:184:THR:HG21	1:B:188:ASN:CB	2.09	0.81
1:B:330:VAL:HG22	1:B:347:THR:HG22	1.63	0.81
1:C:181:LEU:HD21	1:C:247:LYS:HE3	1.62	0.81
1:C:127:LEU:HD21	1:C:236:ILE:HD13	1.63	0.79
1:B:305:ILE:HD11	1:B:336:MSE:HG2	1.65	0.78
1:B:184:THR:CG2	1:B:188:ASN:HB2	2.14	0.78
1:B:356:HIS:ND1	1:B:361:HIS:HE1	1.81	0.78
1:B:325:ILE:HG21	1:B:348:ILE:HD12	1.66	0.77
1:A:135:THR:HG23	2:A:375:HOH:O	1.85	0.76
1:B:87:GLN:HE22	1:B:138:LEU:H	1.30	0.76
1:B:184:THR:HG21	1:B:188:ASN:ND2	1.99	0.76
1:C:161:PRO:HB3	1:C:199:GLY:HA3	1.68	0.76
1:C:173:ILE:HD11	1:C:213:MSE:HE3	1.67	0.75
1:A:117:LEU:HD23	2:A:421:HOH:O	1.87	0.75
1:B:124:LEU:HD21	1:B:183:PRO:HB3	1.68	0.75
1:A:306:SER:OG	1:A:334:VAL:HG23	1.87	0.74
1:A:213:MSE:SE	2:A:442:HOH:O	2.53	0.74
1:A:160:ASN:ND2	1:A:166:GLN:HG3	2.04	0.73
1:A:134:ALA:HB2	1:A:138:LEU:HD11	1.71	0.73
1:B:359:HIS:HD1	1:B:361:HIS:CD2	2.06	0.73
1:B:356:HIS:HD1	1:B:361:HIS:CE1	2.05	0.72
1:A:288:SER:N	1:A:289:PRO:HD3	2.04	0.72
1:B:105:ILE:HD11	1:B:113:VAL:CG1	2.20	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:MSE:HE1	1:B:53:ARG:HH12	1.54	0.72
1:B:173:ILE:CD1	1:B:213:MSE:HE3	2.19	0.72
1:A:207:ASN:HB3	1:A:213:MSE:HE2	1.70	0.72
1:A:63:ASN:C	1:A:63:ASN:HD22	1.97	0.71
1:A:42:MSE:HE1	1:B:53:ARG:HH22	1.55	0.71
1:C:161:PRO:HG3	2:C:412:HOH:O	1.91	0.70
1:B:305:ILE:CD1	1:B:336:MSE:HE3	2.22	0.70
1:A:62:TYR:HD2	1:A:79:LEU:HD11	1.56	0.69
1:B:359:HIS:ND1	1:B:361:HIS:NE2	2.32	0.69
1:A:161:PRO:HB2	1:A:163:ASN:OD1	1.92	0.69
1:C:87:GLN:HE22	1:C:138:LEU:H	1.41	0.69
1:B:63:ASN:HD21	1:B:101:ALA:HA	1.58	0.68
1:A:305:ILE:HB	1:A:334:VAL:HG11	1.73	0.68
1:A:304:ILE:HD13	1:A:335:VAL:C	2.19	0.68
1:B:62:TYR:HB2	1:B:105:ILE:HG23	1.75	0.67
1:A:160:ASN:N	1:A:160:ASN:HD22	1.91	0.67
1:A:334:VAL:HG12	1:A:335:VAL:N	2.09	0.67
1:B:124:LEU:HD21	1:B:183:PRO:HA	1.77	0.67
1:B:124:LEU:HD21	1:B:183:PRO:CB	2.24	0.67
1:C:63:ASN:ND2	1:C:63:ASN:C	2.53	0.67
1:C:185:GLY:CA	1:C:221:ASP:HB2	2.11	0.66
1:B:178:ARG:HH11	1:B:178:ARG:HB3	1.60	0.66
1:A:55:ALA:HB1	1:A:166:GLN:HE22	1.60	0.66
1:A:87:GLN:HE21	1:A:138:LEU:H	1.44	0.66
1:C:299:GLN:HB3	1:C:302:ASP:OD1	1.96	0.66
1:B:189:PHE:HE1	1:B:220:PHE:CD2	2.15	0.65
1:B:197:ASN:CG	1:B:198:HIS:H	2.04	0.65
1:C:124:LEU:HD13	1:C:183:PRO:HB3	1.79	0.65
1:A:79:LEU:HD23	1:A:162:TYR:CB	2.27	0.65
1:B:62:TYR:HB2	1:B:105:ILE:CG2	2.27	0.65
1:B:55:ALA:HB1	1:B:166:GLN:NE2	2.12	0.65
1:C:102:ASP:C	1:C:102:ASP:OD2	2.39	0.64
1:B:356:HIS:CG	1:B:361:HIS:CE1	2.83	0.64
1:B:261:ILE:HG22	1:B:263:GLY:H	1.62	0.64
1:B:356:HIS:HB3	1:B:361:HIS:CE1	2.32	0.64
1:C:258:TYR:HB3	1:C:349:GLN:O	1.98	0.64
1:B:190:LEU:C	1:B:190:LEU:HD23	2.22	0.64
1:A:77:ARG:CZ	2:A:384:HOH:O	2.46	0.63
1:A:259:ILE:HA	1:A:292:PRO:CG	2.28	0.63
1:B:305:ILE:HD11	1:B:336:MSE:HE3	1.79	0.63
1:A:155:VAL:HG13	1:A:205:LEU:CD2	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:308:ASP:O	1:C:309:ASN:CB	2.45	0.62
1:B:258:TYR:CZ	1:B:260:GLY:HA2	2.34	0.62
1:A:147:ARG:HD2	1:A:211:GLU:OE2	2.00	0.62
1:B:262:GLY:O	1:B:285:ASN:HB2	2.00	0.62
1:B:124:LEU:HD21	1:B:183:PRO:CA	2.30	0.62
1:A:127:LEU:HD21	1:A:236:ILE:HD13	1.81	0.61
1:C:327:PRO:HB3	1:C:350:GLU:H	1.65	0.61
1:C:291:GLY:C	1:C:293:ALA:H	2.08	0.61
1:B:124:LEU:CD2	1:B:183:PRO:HB3	2.31	0.61
1:A:125:THR:HG21	1:A:244:ILE:CD1	2.31	0.61
1:A:259:ILE:HA	1:A:292:PRO:HG3	1.81	0.61
1:A:300:VAL:HG22	1:A:300:VAL:O	2.01	0.61
1:C:161:PRO:HB3	1:C:199:GLY:CA	2.30	0.60
1:A:207:ASN:HD22	1:A:213:MSE:HE2	1.64	0.60
1:A:289:PRO:O	1:A:290:ASP:HB3	2.00	0.60
1:A:124:LEU:HD13	1:A:183:PRO:HB3	1.82	0.60
1:A:155:VAL:HG13	1:A:205:LEU:HD22	1.82	0.59
1:B:359:HIS:O	1:B:360:HIS:HB3	2.02	0.59
1:C:158:ILE:HG12	1:C:168:ILE:HD13	1.84	0.59
1:C:351:TYR:CG	1:C:352:PRO:HD2	2.38	0.59
1:C:173:ILE:CD1	1:C:213:MSE:HE3	2.33	0.59
1:C:344:LEU:O	1:C:345:GLN:C	2.45	0.58
1:A:55:ALA:HB1	1:A:166:GLN:NE2	2.17	0.58
1:B:63:ASN:HD22	1:B:64:ARG:N	2.01	0.58
1:B:126:ASP:HB2	1:B:220:PHE:CZ	2.38	0.58
1:C:159:GLY:HA3	1:C:200:ASN:HA	1.85	0.58
1:B:94:ASN:OD1	1:B:202:GLY:N	2.37	0.58
1:B:126:ASP:HB2	1:B:220:PHE:CE2	2.38	0.58
1:C:160:ASN:ND2	1:C:166:GLN:HG3	2.17	0.58
1:B:181:LEU:HD11	1:B:247:LYS:CE	2.26	0.58
1:C:247:LYS:HD3	2:C:433:HOH:O	2.04	0.58
1:C:325:ILE:CD1	1:C:348:ILE:HG13	2.33	0.58
1:A:79:LEU:HD23	1:A:162:TYR:HB3	1.84	0.58
1:B:184:THR:CG2	1:B:188:ASN:ND2	2.62	0.57
1:C:299:GLN:HG3	1:C:300:VAL:N	2.20	0.57
1:A:159:GLY:C	1:A:160:ASN:HD22	2.12	0.57
1:A:236:ILE:HD12	1:A:236:ILE:N	2.20	0.57
1:C:299:GLN:HG3	1:C:300:VAL:H	1.70	0.57
1:B:178:ARG:HB3	1:B:178:ARG:NH1	2.19	0.57
1:B:304:ILE:HD12	1:B:304:ILE:N	2.19	0.57
1:B:85:MSE:HG3	1:B:92:ILE:HG13	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:VAL:HG13	1:C:293:ALA:HB3	1.86	0.56
1:A:239:GLN:H	1:A:239:GLN:NE2	1.91	0.56
1:B:161:PRO:HB2	1:B:163:ASN:OD1	2.05	0.56
1:B:359:HIS:ND1	1:B:361:HIS:CD2	2.74	0.56
1:C:124:LEU:HD13	1:C:183:PRO:CB	2.35	0.56
1:A:42:MSE:HE1	1:B:53:ARG:NH1	2.21	0.56
1:B:297:GLY:HA3	1:B:344:LEU:HD13	1.87	0.56
1:B:306:SER:HA	1:B:312:ALA:HB2	1.86	0.56
1:A:296:ALA:CB	1:A:346:VAL:HG22	2.36	0.56
1:C:239:GLN:H	1:C:239:GLN:NE2	1.95	0.56
1:C:287:VAL:HG12	1:C:287:VAL:O	2.07	0.56
1:A:48:ASN:OD1	1:A:52:ARG:HD3	2.06	0.55
1:B:63:ASN:C	1:B:63:ASN:ND2	2.64	0.55
1:B:259:ILE:C	1:B:292:PRO:HG2	2.31	0.55
1:B:163:ASN:O	1:B:164:LEU:HD22	2.07	0.55
1:B:190:LEU:HD23	1:B:191:GLN:N	2.21	0.55
1:A:350:GLU:O	1:A:351:TYR:C	2.49	0.55
1:C:161:PRO:HD3	1:C:199:GLY:O	2.07	0.55
1:C:284:VAL:HG12	1:C:285:ASN:N	2.20	0.55
1:A:305:ILE:HB	1:A:334:VAL:CG1	2.37	0.55
1:A:42:MSE:HE1	1:B:53:ARG:NH2	2.21	0.55
1:C:127:LEU:CD2	1:C:236:ILE:HD13	2.34	0.54
1:A:63:ASN:CG	1:A:101:ALA:HB2	2.33	0.54
1:B:160:ASN:ND2	1:B:166:GLN:HB2	2.23	0.54
1:B:66:LEU:N	1:B:102:ASP:OD2	2.40	0.54
1:C:239:GLN:HE21	1:C:239:GLN:N	1.96	0.54
1:B:63:ASN:HD22	1:B:63:ASN:C	2.14	0.54
1:B:160:ASN:HD21	1:B:166:GLN:HB2	1.71	0.54
1:B:161:PRO:HB3	1:B:199:GLY:HA2	1.90	0.54
1:A:189:PHE:HE1	1:A:220:PHE:CD2	2.26	0.54
1:C:330:VAL:HG22	1:C:347:THR:CG2	2.17	0.53
1:A:190:LEU:HD23	1:A:190:LEU:C	2.34	0.53
1:C:190:LEU:C	1:C:190:LEU:HD23	2.34	0.53
1:B:79:LEU:CD2	1:B:162:TYR:HB3	2.39	0.53
1:B:109:GLN:HG2	2:B:459:HOH:O	2.07	0.53
1:B:315:ALA:O	1:B:319:MSE:HG2	2.09	0.53
1:C:147:ARG:CD	1:C:211:GLU:OE2	2.57	0.53
1:B:313:ILE:N	1:B:313:ILE:HD12	2.24	0.53
1:C:160:ASN:N	1:C:160:ASN:HD22	2.06	0.53
1:A:288:SER:N	1:A:289:PRO:CD	2.71	0.52
1:B:325:ILE:CG2	1:B:348:ILE:HD12	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:64:ARG:CZ	1:C:103:GLN:OE1	2.57	0.52
1:B:346:VAL:HA	2:B:420:HOH:O	2.08	0.52
1:B:304:ILE:HD12	1:B:304:ILE:H	1.75	0.52
1:B:191:GLN:NE2	1:B:232:ILE:HD11	2.24	0.52
1:A:160:ASN:HA	1:A:166:GLN:HA	1.91	0.52
1:A:77:ARG:NE	2:A:384:HOH:O	2.42	0.52
1:A:318:THR:HA	1:A:321:GLN:OE1	2.10	0.52
1:A:261:ILE:HG22	1:A:262:GLY:N	2.24	0.52
1:B:191:GLN:HE21	1:B:232:ILE:CD1	2.21	0.51
1:B:63:ASN:ND2	1:B:101:ALA:HA	2.25	0.51
1:B:90:TYR:CE2	1:B:131:LYS:HD3	2.46	0.51
1:B:191:GLN:NE2	1:B:232:ILE:CD1	2.73	0.51
1:A:306:SER:H	1:A:334:VAL:CG2	2.23	0.51
1:A:196:ILE:HA	1:A:200:ASN:HD22	1.76	0.51
1:B:181:LEU:CD1	1:B:247:LYS:HE3	2.28	0.51
1:B:356:HIS:CB	1:B:361:HIS:CE1	2.94	0.51
1:C:147:ARG:HD3	1:C:211:GLU:OE2	2.10	0.51
1:C:303:LEU:HD11	1:C:305:ILE:HD11	1.93	0.51
1:B:213:MSE:HE1	2:B:365:HOH:O	2.11	0.51
1:C:346:VAL:HG12	1:C:347:THR:N	2.25	0.51
1:B:361:HIS:OXT	1:B:361:HIS:ND1	2.36	0.50
1:C:230:GLU:O	1:C:230:GLU:HG2	2.11	0.50
1:A:63:ASN:HD22	1:A:64:ARG:N	2.08	0.50
1:A:236:ILE:HD12	1:A:236:ILE:H	1.76	0.50
1:A:305:ILE:HB	1:A:334:VAL:CB	2.42	0.50
1:B:184:THR:HG21	1:B:188:ASN:CG	2.37	0.50
1:B:336:MSE:HE2	1:B:341:GLN:CB	2.42	0.50
1:B:287:VAL:HG12	1:B:287:VAL:O	2.12	0.49
1:C:258:TYR:HB2	1:C:351:TYR:HA	1.94	0.49
1:A:59:VAL:HG12	1:A:108:LEU:HD22	1.94	0.49
1:C:159:GLY:C	1:C:160:ASN:HD22	2.20	0.49
1:A:199:GLY:C	2:A:397:HOH:O	2.55	0.49
1:A:66:LEU:O	1:A:72:ASN:HA	2.13	0.49
1:A:67:ASN:O	1:A:67:ASN:CG	2.55	0.49
1:B:105:ILE:HD11	1:B:113:VAL:HG11	1.95	0.49
1:C:79:LEU:CD2	1:C:162:TYR:HB3	2.43	0.49
1:A:284:VAL:HG12	1:A:284:VAL:O	2.12	0.49
1:A:309:ASN:O	1:A:310:LYS:HD3	2.12	0.49
1:A:297:GLY:HA3	1:A:344:LEU:CB	2.43	0.49
1:B:220:PHE:O	1:B:221:ASP:CB	2.61	0.49
1:B:288:SER:O	1:B:294:ALA:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:ILE:HD12	1:B:334:VAL:CG1	2.35	0.49
1:B:159:GLY:HA3	1:B:200:ASN:HA	1.95	0.48
1:A:190:LEU:HD23	1:A:191:GLN:N	2.28	0.48
1:A:148:VAL:HG12	2:A:399:HOH:O	2.13	0.48
1:C:63:ASN:HD22	1:C:64:ARG:N	2.10	0.48
1:C:194:ALA:O	1:C:196:ILE:HG12	2.14	0.48
1:A:248:LEU:O	1:A:251:ASP:O	2.32	0.48
1:B:147:ARG:HD3	1:B:211:GLU:OE2	2.13	0.48
1:A:159:GLY:HA3	1:A:200:ASN:HA	1.96	0.48
1:A:326:ARG:HG2	1:A:326:ARG:HH11	1.78	0.48
1:C:94:ASN:HB2	1:C:97:VAL:HG23	1.94	0.48
1:A:62:TYR:CD2	1:A:79:LEU:HD11	2.43	0.48
1:C:219:SER:O	1:C:220:PHE:HB2	2.13	0.47
1:B:182:ASN:HB3	1:B:320:ASP:OD1	2.14	0.47
1:C:282:ILE:O	1:C:303:LEU:HA	2.14	0.47
1:B:161:PRO:HB3	1:B:199:GLY:CA	2.44	0.47
1:A:178:ARG:HB3	1:A:178:ARG:NH1	2.15	0.47
1:C:291:GLY:C	1:C:293:ALA:N	2.73	0.47
1:A:135:THR:CG2	2:A:375:HOH:O	2.54	0.47
1:C:299:GLN:CG	1:C:300:VAL:H	2.27	0.47
1:A:156:LEU:CD2	1:A:170:GLN:HB3	2.45	0.47
1:C:261:ILE:HG23	1:C:284:VAL:HG13	1.97	0.47
1:B:96:HIS:HA	1:B:99:ASN:HD22	1.81	0.46
1:A:124:LEU:HD13	1:A:183:PRO:CB	2.46	0.46
1:A:134:ALA:HB2	1:A:138:LEU:CD1	2.43	0.46
1:B:55:ALA:HB1	1:B:166:GLN:HE22	1.78	0.46
1:A:42:MSE:CE	1:B:53:ARG:HH12	2.25	0.46
1:A:304:ILE:HG22	1:A:305:ILE:N	2.30	0.46
1:C:322:VAL:O	1:C:325:ILE:HG12	2.15	0.46
1:A:78:THR:HG23	2:A:393:HOH:O	2.16	0.46
1:B:156:LEU:CD2	1:B:170:GLN:HB3	2.45	0.46
1:A:155:VAL:HG13	1:A:205:LEU:HD21	1.98	0.46
1:A:284:VAL:CG2	1:A:304:ILE:HG12	2.46	0.46
1:B:148:VAL:HG22	1:B:150:HIS:CE1	2.51	0.46
1:B:160:ASN:ND2	1:B:166:GLN:CB	2.79	0.46
1:C:160:ASN:ND2	1:C:160:ASN:N	2.64	0.46
1:B:197:ASN:CG	1:B:198:HIS:N	2.70	0.46
1:C:330:VAL:CG2	1:C:347:THR:HG22	2.17	0.46
1:B:48:ASN:OD1	1:B:52:ARG:NH1	2.42	0.45
1:A:261:ILE:CG2	1:A:262:GLY:N	2.79	0.45
1:C:230:GLU:O	1:C:230:GLU:CG	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:HIS:CB	1:B:361:HIS:HE1	2.28	0.45
1:A:63:ASN:C	1:A:63:ASN:ND2	2.69	0.45
1:A:126:ASP:HB2	1:A:220:PHE:CZ	2.52	0.45
1:C:48:ASN:OD1	1:C:52:ARG:NH1	2.40	0.45
1:C:219:SER:O	1:C:220:PHE:CB	2.64	0.45
1:A:59:VAL:CG1	1:A:108:LEU:HD22	2.47	0.45
1:A:161:PRO:HD3	1:A:199:GLY:O	2.17	0.45
1:A:290:ASP:CG	1:A:290:ASP:O	2.59	0.45
1:B:305:ILE:CD1	1:B:336:MSE:HG2	2.42	0.45
1:C:181:LEU:CD2	1:C:247:LYS:HE3	2.39	0.45
1:C:287:VAL:CG1	1:C:293:ALA:HB3	2.47	0.45
1:B:191:GLN:HE21	1:B:232:ILE:HD11	1.82	0.45
1:C:109:GLN:NE2	1:C:160:ASN:OD1	2.50	0.45
1:B:65:GLY:HA2	1:B:102:ASP:CG	2.31	0.45
1:B:305:ILE:HD11	1:B:336:MSE:CE	2.47	0.45
1:B:306:SER:HB2	1:B:310:LYS:O	2.17	0.45
1:C:262:GLY:O	1:C:285:ASN:HB3	2.17	0.45
1:A:48:ASN:O	1:A:52:ARG:HG3	2.17	0.44
1:C:161:PRO:CG	2:C:412:HOH:O	2.60	0.44
1:A:126:ASP:HB2	1:A:220:PHE:CE2	2.52	0.44
1:B:213:MSE:CE	2:B:365:HOH:O	2.64	0.44
1:C:303:LEU:O	1:C:304:ILE:HD13	2.17	0.44
1:A:303:LEU:C	1:A:303:LEU:HD23	2.42	0.44
1:B:60:ASN:OD1	1:B:160:ASN:O	2.35	0.44
1:A:326:ARG:HG2	1:A:326:ARG:NH1	2.31	0.44
1:A:218:LEU:HD23	1:A:218:LEU:HA	1.84	0.44
1:A:287:VAL:C	1:A:289:PRO:HD3	2.42	0.44
1:A:331:ILE:HG12	1:A:348:ILE:HD11	1.98	0.44
1:B:183:PRO:HG3	1:B:351:TYR:CZ	2.52	0.44
1:A:85:MSE:HE1	1:A:143:ILE:HD11	1.99	0.44
1:A:220:PHE:O	1:A:221:ASP:CB	2.65	0.44
1:C:330:VAL:HA	1:C:346:VAL:O	2.18	0.44
1:C:185:GLY:O	1:C:186:ARG:C	2.61	0.43
1:A:124:LEU:HD12	1:A:256:ARG:CZ	2.48	0.43
1:A:307:VAL:HG21	1:A:321:GLN:OE1	2.17	0.43
1:C:85:MSE:HG3	1:C:92:ILE:HG13	1.99	0.43
1:B:197:ASN:ND2	1:B:198:HIS:H	2.16	0.43
1:C:282:ILE:CD1	1:C:312:ALA:HB1	2.48	0.43
1:B:258:TYR:OH	1:B:260:GLY:HA2	2.18	0.43
1:C:94:ASN:OD1	1:C:202:GLY:N	2.50	0.43
1:A:160:ASN:ND2	1:A:160:ASN:N	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:ASN:ND2	1:B:164:LEU:HD23	2.34	0.43
1:C:79:LEU:HD23	1:C:162:TYR:HB3	1.99	0.43
1:A:197:ASN:ND2	1:B:230:GLU:HG3	2.34	0.43
1:A:294:ALA:C	1:A:296:ALA:H	2.25	0.43
1:C:58:VAL:O	1:C:160:ASN:OD1	2.35	0.43
1:A:321:GLN:N	2:A:470:HOH:O	2.33	0.43
1:B:259:ILE:O	1:B:292:PRO:HG2	2.18	0.43
1:A:284:VAL:HG23	1:A:304:ILE:HG12	2.00	0.43
1:C:291:GLY:O	1:C:293:ALA:N	2.52	0.43
1:B:182:ASN:OD1	1:B:183:PRO:HD2	2.19	0.43
1:B:184:THR:HG23	1:B:188:ASN:HD22	1.78	0.43
1:B:300:VAL:HG13	1:B:301:ASN:N	2.34	0.43
1:A:90:TYR:CE2	1:A:131:LYS:HD3	2.53	0.42
1:A:296:ALA:HB3	1:A:346:VAL:HG22	2.01	0.42
1:B:47:TYR:HB3	1:B:156:LEU:HD11	2.01	0.42
1:C:179:ILE:HG21	1:C:239:GLN:HB2	2.00	0.42
1:C:194:ALA:O	1:C:195:SER:C	2.62	0.42
1:B:196:ILE:HG12	1:B:233:GLY:HA3	2.01	0.42
1:C:213:MSE:CE	2:C:365:HOH:O	2.67	0.42
1:A:245:MSE:SE	2:A:478:HOH:O	2.87	0.42
1:B:291:GLY:H	1:B:294:ALA:HB3	1.83	0.42
1:A:195:SER:O	1:A:200:ASN:ND2	2.53	0.42
1:A:288:SER:C	1:A:290:ASP:H	2.27	0.42
1:B:147:ARG:CD	1:B:211:GLU:OE2	2.67	0.42
1:B:313:ILE:HD12	1:B:313:ILE:H	1.83	0.42
1:C:284:VAL:CG1	1:C:285:ASN:N	2.83	0.42
1:A:85:MSE:HG3	1:A:92:ILE:HG13	2.01	0.42
1:A:98:ILE:HD12	1:A:98:ILE:C	2.45	0.42
1:B:220:PHE:O	1:B:221:ASP:HB3	2.20	0.42
1:A:79:LEU:HD23	1:A:162:TYR:HB2	1.99	0.42
1:A:167:THR:HG22	2:B:433:HOH:O	2.20	0.42
1:A:259:ILE:HA	1:A:292:PRO:HG2	2.00	0.42
1:A:286:GLU:OE1	1:A:286:GLU:C	2.63	0.42
1:A:184:THR:OG1	1:A:188:ASN:ND2	2.48	0.41
1:B:239:GLN:H	1:B:239:GLN:NE2	1.98	0.41
1:A:320:ASP:HB2	2:A:470:HOH:O	2.19	0.41
1:C:196:ILE:HA	1:C:200:ASN:HD22	1.85	0.41
1:C:284:VAL:HG23	1:C:302:ASP:O	2.20	0.41
1:B:185:GLY:HA2	1:B:358:HIS:CE1	2.55	0.41
1:C:147:ARG:HD2	1:C:211:GLU:OE2	2.20	0.41
1:A:186:ARG:CB	1:A:221:ASP:HB2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:LEU:HD23	1:C:191:GLN:N	2.35	0.41
1:A:321:GLN:O	1:A:325:ILE:HG12	2.21	0.41
1:B:299:GLN:OE1	1:B:299:GLN:HA	2.19	0.41
1:B:284:VAL:HB	1:B:300:VAL:O	2.20	0.41
1:A:42:MSE:CE	1:B:53:ARG:HH22	2.30	0.41
1:B:184:THR:HG22	1:B:185:GLY:O	2.21	0.41
1:B:220:PHE:HE1	1:B:358:HIS:CD2	2.39	0.41
1:C:179:ILE:CG2	1:C:239:GLN:HB2	2.51	0.41
1:C:283:VAL:HG23	1:C:283:VAL:O	2.20	0.41
1:C:315:ALA:O	1:C:319:MSE:HG2	2.21	0.41
1:A:62:TYR:HD2	1:A:79:LEU:CD1	2.30	0.41
1:A:288:SER:C	1:A:290:ASP:N	2.78	0.41
1:A:304:ILE:HD12	1:A:333:VAL:CG1	2.51	0.41
1:B:42:MSE:HA	1:B:42:MSE:HE2	2.02	0.41
1:B:229:PRO:HA	1:B:232:ILE:HG23	2.02	0.41
1:B:230:GLU:CD	1:B:230:GLU:N	2.78	0.41
1:A:317:GLU:C	2:A:470:HOH:O	2.64	0.41
1:C:95:LYS:HE3	2:C:407:HOH:O	2.21	0.41
1:B:335:VAL:HG12	1:B:336:MSE:N	2.36	0.40
1:C:196:ILE:HD12	1:C:216:ASN:HB3	2.03	0.40
1:A:99:ASN:ND2	2:A:481:HOH:O	2.46	0.40
1:A:330:VAL:CG2	1:A:347:THR:HG22	2.35	0.40
1:B:105:ILE:HD11	1:B:113:VAL:HG13	1.98	0.40
1:B:218:LEU:HD23	1:B:218:LEU:HA	1.91	0.40
1:C:325:ILE:HD12	1:C:348:ILE:HG13	2.03	0.40
1:C:318:THR:O	1:C:322:VAL:HG23	2.22	0.40
1:B:59:VAL:HG11	1:B:108:LEU:HD22	2.03	0.40
1:B:190:LEU:C	1:B:190:LEU:CD2	2.93	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ASN:ND2	1:B:361:HIS:NE2[1_554]	2.17	0.03

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	275/320 (86%)	244 (89%)	22 (8%)	9 (3%)	3	2
1	B	286/320 (89%)	253 (88%)	25 (9%)	8 (3%)	4	2
1	C	273/320 (85%)	228 (84%)	33 (12%)	12 (4%)	2	1
All	All	834/960 (87%)	725 (87%)	80 (10%)	29 (4%)	3	1

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	219	SER
1	A	313	ILE
1	A	334	VAL
1	B	219	SER
1	B	300	VAL
1	C	220	PHE
1	C	343	THR
1	A	343	THR
1	B	186	ARG
1	B	298	ILE
1	C	313	ILE
1	C	345	GLN
1	B	220	PHE
1	C	299	GLN
1	A	301	ASN
1	B	74	LEU
1	C	103	GLN
1	A	351	TYR
1	C	74	LEU
1	C	219	SER
1	C	287	VAL
1	C	292	PRO
1	A	311	PRO
1	C	300	VAL

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Mol	Chain	Res	Type
1	A	298	ILE
1	B	313	ILE
1	A	284	VAL
1	B	297	GLY
1	C	284	VAL

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	217/253 (86%)	200 (92%)	17 (8%)	11	16
1	B	232/253 (92%)	216 (93%)	16 (7%)	14	20
1	C	214/253 (85%)	199 (93%)	15 (7%)	14	19
All	All	663/759 (87%)	615 (93%)	48 (7%)	13	18

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	79	LEU
1	A	108	LEU
1	A	118	LEU
1	A	119	VAL
1	A	147	ARG
1	A	155	VAL
1	A	160	ASN
1	A	168	ILE
1	A	178	ARG
1	A	201	SER
1	A	205	LEU
1	A	219	SER
1	A	220	PHE
1	A	239	GLN
1	A	240	LEU
1	A	286	GLU

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Mol	Chain	Res	Type
1	B	106	VAL
1	B	108	LEU
1	B	118	LEU
1	B	119	VAL
1	B	148	VAL
1	B	160	ASN
1	B	164	LEU
1	B	184	THR
1	B	201	SER
1	B	205	LEU
1	B	220	PHE
1	B	230	GLU
1	B	232	ILE
1	B	239	GLN
1	B	240	LEU
1	B	361	HIS
1	C	63	ASN
1	C	102	ASP
1	C	108	LEU
1	C	118	LEU
1	C	119	VAL
1	C	140	THR
1	C	168	ILE
1	C	187	GLN
1	C	197	ASN
1	C	205	LEU
1	C	220	PHE
1	C	232	ILE
1	C	239	GLN
1	C	240	LEU
1	C	286	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	71	HIS
1	A	73	GLN
1	A	87	GLN
1	A	109	GLN
1	A	133	ASN
1	A	160	ASN

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Mol	Chain	Res	Type
1	A	166	GLN
1	A	170	GLN
1	A	182	ASN
1	A	188	ASN
1	A	239	GLN
1	A	299	GLN
1	A	309	ASN
1	A	345	GLN
1	B	60	ASN
1	B	63	ASN
1	B	87	GLN
1	B	109	GLN
1	B	150	HIS
1	B	160	ASN
1	B	166	GLN
1	B	170	GLN
1	B	188	ASN
1	B	191	GLN
1	B	239	GLN
1	B	349	GLN
1	B	358	HIS
1	C	63	ASN
1	C	87	GLN
1	C	96	HIS
1	C	109	GLN
1	C	160	ASN
1	C	166	GLN
1	C	170	GLN
1	C	188	ASN
1	C	239	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 [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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	77:ARG	C	78:THR	N	1.11
1	B	77:ARG	C	78:THR	N	0.93

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	278/320 (86%)	0.39	36 (12%) 7 8	14, 44, 139, 145	0
1	B	277/320 (86%)	0.60	48 (17%) 4 4	24, 54, 124, 143	0
1	C	265/320 (82%)	0.57	41 (15%) 5 6	24, 54, 142, 149	0
All	All	820/960 (85%)	0.52	125 (15%) 5 6	14, 51, 137, 149	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	66	LEU	7.5
1	C	66	LEU	5.9
1	A	220	PHE	5.7
1	B	229	PRO	5.4
1	A	221	ASP	5.3
1	B	354	THR	5.1
1	B	360	HIS	4.8
1	C	160	ASN	4.8
1	B	298	ILE	4.5
1	B	220	PHE	4.4
1	A	161	PRO	4.4
1	C	309	ASN	4.3
1	C	264	ARG	4.1
1	C	220	PHE	4.1
1	A	307	VAL	4.1
1	B	67	ASN	4.1
1	A	162	TYR	4.0
1	A	341	GLN	3.9
1	C	65	GLY	3.9
1	B	65	GLY	3.8
1	B	304	ILE	3.7
1	B	300	VAL	3.7
1	C	67	ASN	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	162	TYR	3.6
1	A	287	VAL	3.6
1	C	307	VAL	3.6
1	B	312	ALA	3.6
1	B	305	ILE	3.5
1	B	101	ALA	3.4
1	B	287	VAL	3.4
1	B	165	GLY	3.4
1	A	293	ALA	3.3
1	C	199	GLY	3.3
1	C	300	VAL	3.3
1	A	316	LEU	3.3
1	A	304	ILE	3.3
1	B	361	HIS	3.3
1	C	287	VAL	3.3
1	A	160	ASN	3.2
1	A	334	VAL	3.2
1	C	162	TYR	3.2
1	C	161	PRO	3.2
1	C	342	LEU	3.2
1	C	313	ILE	3.2
1	C	343	THR	3.1
1	B	295	ASN	3.1
1	C	305	ILE	3.1
1	B	307	VAL	3.1
1	A	305	ILE	3.0
1	B	160	ASN	3.0
1	A	294	ALA	3.0
1	B	342	LEU	3.0
1	B	197	ASN	3.0
1	B	297	GLY	3.0
1	A	342	LEU	3.0
1	B	334	VAL	3.0
1	B	219	SER	3.0
1	C	294	ALA	2.9
1	C	164	LEU	2.9
1	A	313	ILE	2.9
1	A	315	ALA	2.8
1	C	344	LEU	2.8
1	B	356	HIS	2.8
1	B	341	GLN	2.7
1	B	221	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	221	ASP	2.7
1	C	218	LEU	2.7
1	A	343	THR	2.7
1	C	331	ILE	2.7
1	C	332	PRO	2.7
1	C	330	VAL	2.6
1	A	297	GLY	2.6
1	B	290	ASP	2.6
1	C	292	PRO	2.6
1	A	300	VAL	2.6
1	C	289	PRO	2.6
1	B	344	LEU	2.6
1	C	101	ALA	2.5
1	A	292	PRO	2.5
1	B	335	VAL	2.5
1	A	309	ASN	2.5
1	B	296	ALA	2.5
1	B	260	GLY	2.5
1	A	306	SER	2.4
1	B	283	VAL	2.4
1	C	301	ASN	2.4
1	B	264	ARG	2.4
1	C	263	GLY	2.4
1	A	298	ILE	2.4
1	C	298	ILE	2.4
1	B	355	ASN	2.3
1	B	303	LEU	2.3
1	B	161	PRO	2.3
1	B	294	ALA	2.3
1	A	219	SER	2.3
1	A	264	ARG	2.3
1	C	288	SER	2.3
1	C	303	LEU	2.3
1	B	263	GLY	2.3
1	C	299	GLN	2.2
1	C	261	ILE	2.2
1	C	312	ALA	2.2
1	A	303	LEU	2.2
1	A	163	ASN	2.2
1	A	134	ALA	2.2
1	A	348	ILE	2.2
1	C	163	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	218	LEU	2.1
1	A	333	VAL	2.1
1	B	299	GLN	2.1
1	B	259	ILE	2.1
1	C	327	PRO	2.1
1	A	346	VAL	2.1
1	A	345	GLN	2.0
1	B	292	PRO	2.0
1	B	311	PRO	2.0
1	B	284	VAL	2.0
1	C	219	SER	2.0
1	B	353	ALA	2.0
1	C	100	ASP	2.0
1	A	311	PRO	2.0
1	B	289	PRO	2.0
1	B	348	ILE	2.0
1	C	311	PRO	2.0
1	A	310	LYS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.