



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:00 PM UTC

PDB ID : 2ZNJ / pdb_00002znj
Title : Crystal structure of Pyrrolsyt-tRNA synthetase from *Desulfitobacterium hafniense*
Authors : Nozawa, K.; Araiso, Y.; Soll, D.; Ishitani, R.; Nureki, O.
Deposited on : 2008-04-25
Resolution : 2.50 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

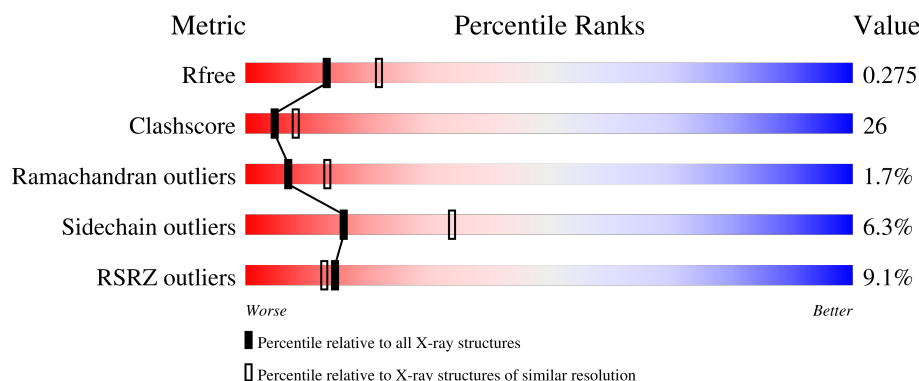
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	308	12% 51% 37% • 10%
1	B	308	9% 55% 29% 6% • 9%
1	C	308	4% 56% 30% • 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7189 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 Putative uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			2255	1434	396	413	12			
1	B	279	Total	C	N	O	S	0	0	0
			2263	1440	397	414	12			
1	C	278	Total	C	N	O	S	0	0	0
			2255	1434	396	413	12			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP B0S4P3
A	-18	GLY	-	expression tag	UNP B0S4P3
A	-17	SER	-	expression tag	UNP B0S4P3
A	-16	SER	-	expression tag	UNP B0S4P3
A	-15	HIS	-	expression tag	UNP B0S4P3
A	-14	HIS	-	expression tag	UNP B0S4P3
A	-13	HIS	-	expression tag	UNP B0S4P3
A	-12	HIS	-	expression tag	UNP B0S4P3
A	-11	HIS	-	expression tag	UNP B0S4P3
A	-10	HIS	-	expression tag	UNP B0S4P3
A	-9	SER	-	expression tag	UNP B0S4P3
A	-8	SER	-	expression tag	UNP B0S4P3
A	-7	GLY	-	expression tag	UNP B0S4P3
A	-6	LEU	-	expression tag	UNP B0S4P3
A	-5	VAL	-	expression tag	UNP B0S4P3
A	-4	PRO	-	expression tag	UNP B0S4P3
A	-3	ARG	-	expression tag	UNP B0S4P3
A	-2	GLY	-	expression tag	UNP B0S4P3
A	-1	SER	-	expression tag	UNP B0S4P3
A	0	HIS	-	expression tag	UNP B0S4P3
B	-19	MET	-	expression tag	UNP B0S4P3
B	-18	GLY	-	expression tag	UNP B0S4P3
B	-17	SER	-	expression tag	UNP B0S4P3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	expression tag	UNP B0S4P3
B	-15	HIS	-	expression tag	UNP B0S4P3
B	-14	HIS	-	expression tag	UNP B0S4P3
B	-13	HIS	-	expression tag	UNP B0S4P3
B	-12	HIS	-	expression tag	UNP B0S4P3
B	-11	HIS	-	expression tag	UNP B0S4P3
B	-10	HIS	-	expression tag	UNP B0S4P3
B	-9	SER	-	expression tag	UNP B0S4P3
B	-8	SER	-	expression tag	UNP B0S4P3
B	-7	GLY	-	expression tag	UNP B0S4P3
B	-6	LEU	-	expression tag	UNP B0S4P3
B	-5	VAL	-	expression tag	UNP B0S4P3
B	-4	PRO	-	expression tag	UNP B0S4P3
B	-3	ARG	-	expression tag	UNP B0S4P3
B	-2	GLY	-	expression tag	UNP B0S4P3
B	-1	SER	-	expression tag	UNP B0S4P3
B	0	HIS	-	expression tag	UNP B0S4P3
C	-19	MET	-	expression tag	UNP B0S4P3
C	-18	GLY	-	expression tag	UNP B0S4P3
C	-17	SER	-	expression tag	UNP B0S4P3
C	-16	SER	-	expression tag	UNP B0S4P3
C	-15	HIS	-	expression tag	UNP B0S4P3
C	-14	HIS	-	expression tag	UNP B0S4P3
C	-13	HIS	-	expression tag	UNP B0S4P3
C	-12	HIS	-	expression tag	UNP B0S4P3
C	-11	HIS	-	expression tag	UNP B0S4P3
C	-10	HIS	-	expression tag	UNP B0S4P3
C	-9	SER	-	expression tag	UNP B0S4P3
C	-8	SER	-	expression tag	UNP B0S4P3
C	-7	GLY	-	expression tag	UNP B0S4P3
C	-6	LEU	-	expression tag	UNP B0S4P3
C	-5	VAL	-	expression tag	UNP B0S4P3
C	-4	PRO	-	expression tag	UNP B0S4P3
C	-3	ARG	-	expression tag	UNP B0S4P3
C	-2	GLY	-	expression tag	UNP B0S4P3
C	-1	SER	-	expression tag	UNP B0S4P3
C	0	HIS	-	expression tag	UNP B0S4P3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	125	Total O 125 125	0	0

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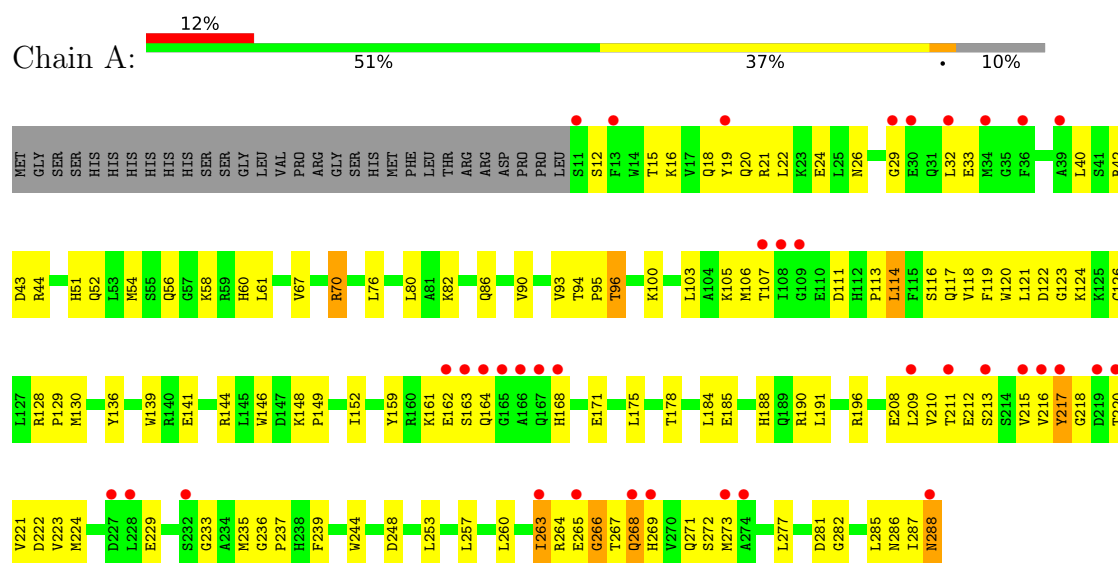
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	123	Total 123	O 123	0	0
2	C	168	Total 168	O 168	0	0

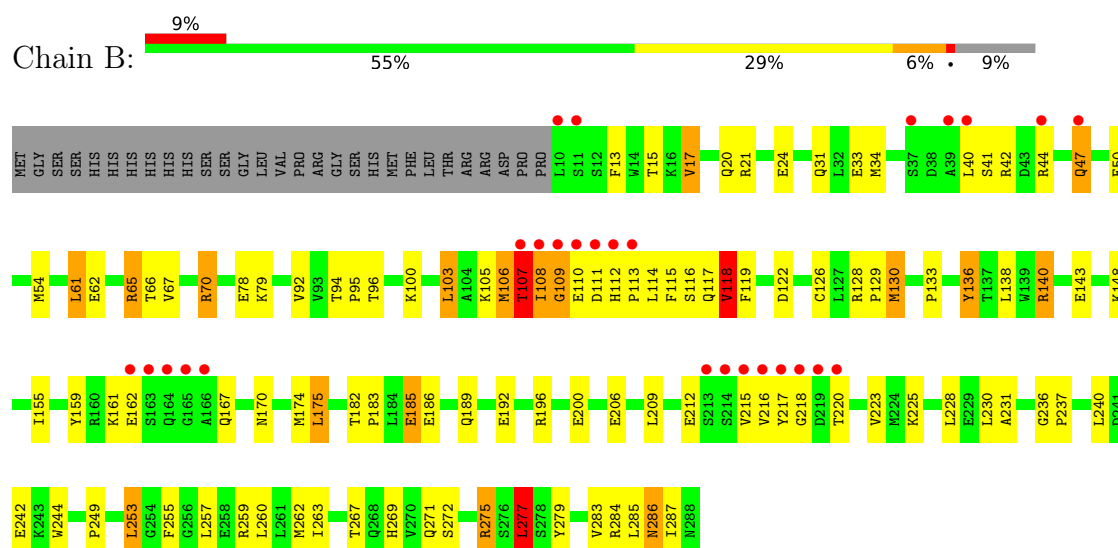
3 Residue-property plots

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: Putative uncharacterized protein

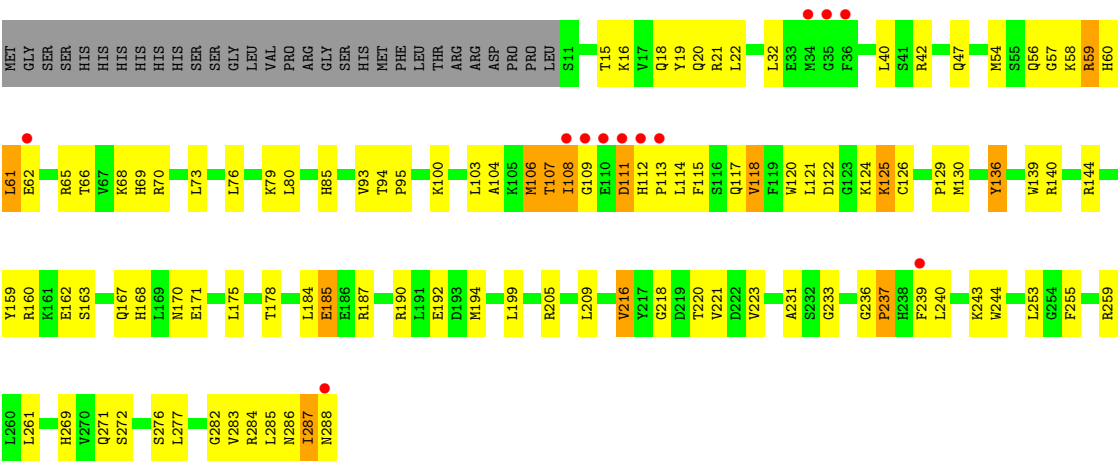


• Molecule 1: Putative uncharacterized protein



• Molecule 1: Putative uncharacterized protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	124.92Å 128.56Å 124.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.67 – 2.50 44.67 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.67-2.50) 100.0 (44.67-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.08Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.214 , 0.275 0.214 , 0.275	Depositor DCC
R_{free} test set	1768 reflections (2.98%)	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.571	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7189	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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	0.42	0/2303	0.92	9/3105 (0.3%)
1	B	0.42	0/2311	0.94	8/3116 (0.3%)
1	C	0.43	0/2303	0.96	12/3105 (0.4%)
All	All	0.42	0/6917	0.94	29/9326 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	MET	N-CA-C	-9.26	103.15	112.97
1	A	123	GLY	N-CA-C	-8.68	103.94	115.32
1	B	110	GLU	N-CA-C	-7.51	104.09	113.18
1	C	216	VAL	N-CA-C	7.34	118.28	111.45
1	C	277	LEU	N-CA-C	-7.13	105.12	114.31
1	A	285	LEU	N-CA-C	7.12	121.20	112.23
1	C	106	MET	N-CA-C	-6.83	104.93	112.72
1	C	287	ILE	N-CA-C	6.59	119.42	109.20
1	C	285	LEU	N-CA-C	6.43	120.95	113.18
1	A	277	LEU	N-CA-C	-5.87	105.96	114.12
1	B	118	VAL	N-CA-C	5.68	117.78	108.97
1	C	118	VAL	N-CA-C	5.68	117.78	108.97
1	B	257	LEU	N-CA-C	5.68	117.47	111.28
1	C	121	LEU	N-CA-C	-5.66	106.51	113.41
1	C	125	LYS	N-CA-C	5.64	118.74	109.72
1	B	109	GLY	N-CA-C	-5.63	99.83	113.18
1	A	129	PRO	N-CA-C	-5.60	106.80	114.98
1	A	164	GLN	N-CA-C	-5.60	105.05	113.89
1	B	170	ASN	N-CA-C	-5.60	105.18	112.23
1	C	66	THR	N-CA-C	5.49	120.25	113.50
1	A	257	LEU	N-CA-C	5.47	117.24	111.28
1	A	96	THR	N-CA-C	-5.45	106.14	112.89
1	B	275	ARG	N-CA-C	-5.39	102.07	110.10
1	C	57	GLY	N-CA-C	-5.17	106.15	112.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	276	SER	N-CA-C	5.15	116.12	108.60
1	B	277	LEU	N-CA-C	-5.12	107.70	114.31
1	C	85	HIS	N-CA-C	-5.10	105.72	111.28
1	A	121	LEU	N-CA-C	-5.10	106.75	113.12
1	A	287	ILE	N-CA-C	5.05	116.13	108.96

There are no chirality outliers.

There are no planarity outliers.

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	2255	0	2249	102	0
1	B	2263	0	2260	142	0
1	C	2255	0	2249	105	0
2	A	125	0	0	13	0
2	B	123	0	0	7	0
2	C	168	0	0	13	0
All	All	7189	0	6758	345	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 26.

All (345) 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:C:104:ALA:HA	1:C:108:ILE:HB	1.32	1.10
1:B:108:ILE:HG12	1:B:109:GLY:H	1.00	1.06
1:C:107:THR:HG23	1:C:240:LEU:HD21	1.33	1.06
1:C:54:MET:HE2	1:C:287:ILE:HA	1.39	1.00
1:C:108:ILE:HG12	1:C:112:HIS:HB3	1.44	1.00
1:C:108:ILE:HD11	1:C:114:LEU:HB3	1.46	0.97
1:B:140:ARG:HH11	1:B:140:ARG:HB3	1.29	0.96
1:B:31:GLN:HA	1:B:34:MET:HE3	1.46	0.94
1:B:108:ILE:HG12	1:B:109:GLY:N	1.81	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:VAL:HG13	1:A:217:TYR:H	1.34	0.92
1:B:103:LEU:HD11	1:B:114:LEU:HD22	1.49	0.92
1:C:21:ARG:HH21	1:C:54:MET:HE3	1.32	0.91
1:C:129:PRO:HG2	1:C:130:MET:HE2	1.53	0.89
1:B:108:ILE:CG1	1:B:109:GLY:H	1.85	0.89
1:B:15:THR:HG23	1:B:42:ARG:HH12	1.39	0.86
1:C:21:ARG:NH2	1:C:54:MET:HE3	1.90	0.85
1:B:228:LEU:HD13	1:B:263:ILE:HD12	1.58	0.84
1:C:112:HIS:ND1	1:C:113:PRO:HD2	1.93	0.84
1:A:52:GLN:HE21	1:A:56:GLN:HE21	1.26	0.83
1:C:58:LYS:HE3	1:C:288:ASN:HB2	1.62	0.81
1:B:113:PRO:O	1:B:117:GLN:HG3	1.82	0.79
1:C:54:MET:HE2	1:C:287:ILE:CA	2.12	0.78
1:C:107:THR:CG2	1:C:240:LEU:HD21	2.12	0.78
1:C:100:LYS:HD3	2:C:433:HOH:O	1.83	0.77
1:B:15:THR:HG23	1:B:42:ARG:NH1	1.99	0.77
1:B:112:HIS:ND1	1:B:113:PRO:HD2	1.99	0.76
1:A:58:LYS:HD2	1:A:288:ASN:HB2	1.66	0.76
1:A:233:GLY:HA3	1:A:253:LEU:HD12	1.67	0.75
1:B:108:ILE:CD1	1:B:114:LEU:HB3	2.17	0.75
1:C:112:HIS:HD2	1:C:114:LEU:HB2	1.51	0.75
1:C:118:VAL:CG1	1:C:126:CYS:HB2	2.17	0.74
1:A:269:HIS:ND1	1:A:271:GLN:HG2	2.01	0.74
1:C:112:HIS:HE1	1:C:216:VAL:HG22	1.52	0.73
1:B:209:LEU:CD2	1:B:223:VAL:HG22	2.19	0.73
1:C:108:ILE:HG21	1:C:115:PHE:CE2	2.25	0.72
1:A:162:GLU:HA	2:A:410:HOH:O	1.88	0.71
1:C:112:HIS:CE1	1:C:216:VAL:HG22	2.25	0.71
1:C:61:LEU:CD2	1:C:283:VAL:HG11	2.22	0.70
1:B:108:ILE:HD12	1:B:114:LEU:HB3	1.72	0.70
1:B:269:HIS:HB3	1:B:271:GLN:HE21	1.57	0.69
1:C:112:HIS:CG	1:C:113:PRO:HD2	2.28	0.69
1:B:40:LEU:HB2	2:B:429:HOH:O	1.91	0.69
1:A:118:VAL:HG11	1:A:126:CYS:SG	2.33	0.68
1:B:129:PRO:HG2	1:B:130:MET:HE2	1.76	0.68
1:B:112:HIS:HD2	1:B:114:LEU:HB2	1.59	0.67
1:B:284:ARG:NH1	1:B:286:ASN:HB2	2.09	0.67
1:C:104:ALA:HA	1:C:108:ILE:CB	2.18	0.67
1:C:56:GLN:HA	1:C:59:ARG:HD2	1.75	0.67
1:A:224:MET:HE3	1:A:229:GLU:HB3	1.76	0.66
1:A:26:ASN:OD1	1:A:267:THR:HG23	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:LYS:HE2	1:A:248:ASP:CG	2.20	0.66
1:A:263:ILE:O	1:A:263:ILE:HG22	1.96	0.65
1:C:117:GLN:HE21	1:C:160:ARG:HH12	1.42	0.65
1:C:112:HIS:HB2	2:C:401:HOH:O	1.97	0.65
1:B:108:ILE:HG13	1:B:112:HIS:HB3	1.76	0.65
1:B:259:ARG:O	1:B:263:ILE:HG12	1.96	0.65
1:B:54:MET:HE2	1:B:287:ILE:HG22	1.79	0.65
1:B:114:LEU:HA	1:B:117:GLN:OE1	1.97	0.64
1:C:113:PRO:O	1:C:117:GLN:HG2	1.98	0.64
1:A:213:SER:HA	1:A:218:GLY:O	1.98	0.64
1:C:184:LEU:HA	1:C:187:ARG:HD3	1.80	0.64
1:B:103:LEU:HD12	1:B:115:PHE:CE1	2.32	0.64
1:B:112:HIS:CD2	1:B:114:LEU:HB2	2.32	0.64
1:B:106:MET:HE2	1:B:244:TRP:HE1	1.62	0.64
1:B:117:GLN:HG2	1:B:162:GLU:OE1	1.98	0.64
1:A:52:GLN:HE21	1:A:56:GLN:NE2	1.96	0.64
1:B:47:GLN:HE21	1:B:47:GLN:C	2.05	0.64
1:A:114:LEU:HD21	1:A:130:MET:HE3	1.79	0.63
1:C:175:LEU:C	1:C:175:LEU:HD23	2.23	0.63
1:C:61:LEU:HD21	1:C:283:VAL:HG11	1.80	0.63
1:A:163:SER:HB2	1:A:168:HIS:HB3	1.79	0.63
1:C:106:MET:HE2	1:C:244:TRP:HE1	1.64	0.63
1:A:119:PHE:CZ	1:A:161:LYS:HA	2.34	0.62
1:B:107:THR:HG22	1:B:107:THR:O	1.99	0.62
1:C:218:GLY:HA3	1:C:239:PHE:HE2	1.64	0.62
1:A:95:PRO:HD2	1:B:277:LEU:HD13	1.80	0.62
1:B:54:MET:HE2	1:B:287:ILE:HA	1.81	0.62
1:A:224:MET:HG2	1:A:229:GLU:HA	1.81	0.62
1:A:19:TYR:CE1	1:A:32:LEU:HD11	2.35	0.61
1:C:185:GLU:H	1:C:185:GLU:CD	2.07	0.61
1:B:40:LEU:HD13	1:B:40:LEU:C	2.25	0.61
1:B:62:GLU:CD	1:B:65:ARG:HH12	2.09	0.61
1:B:262:MET:HG3	1:B:267:THR:HG23	1.83	0.60
1:A:148:LYS:HE2	1:A:248:ASP:OD2	2.01	0.60
1:A:273:MET:HB2	2:A:320:HOH:O	2.01	0.60
1:B:140:ARG:HH11	1:B:140:ARG:CB	2.07	0.60
1:C:205:ARG:HG2	2:C:345:HOH:O	2.02	0.60
1:B:62:GLU:CD	1:B:65:ARG:NH1	2.60	0.60
1:A:32:LEU:HD23	1:A:32:LEU:O	2.02	0.59
1:A:236:GLY:HA2	1:A:237:PRO:C	2.27	0.59
1:C:93:VAL:O	1:C:93:VAL:HG13	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:MET:HE2	1:A:244:TRP:HE1	1.67	0.59
1:A:141:GLU:HG2	2:A:325:HOH:O	2.03	0.59
1:C:118:VAL:HG11	1:C:126:CYS:HB2	1.85	0.58
1:A:163:SER:HA	1:A:168:HIS:HD2	1.67	0.58
1:C:236:GLY:HA2	1:C:237:PRO:C	2.28	0.58
1:A:191:LEU:HD13	1:A:221:VAL:HG21	1.85	0.58
1:A:21:ARG:HH21	1:A:54:MET:CE	2.17	0.58
1:B:108:ILE:HD11	1:B:112:HIS:O	2.04	0.57
1:A:114:LEU:HD21	1:A:130:MET:CE	2.32	0.57
1:A:216:VAL:HG13	1:A:217:TYR:N	2.14	0.57
1:B:225:LYS:HB2	1:B:230:LEU:HD11	1.86	0.57
1:C:20:GLN:HG2	2:C:337:HOH:O	2.04	0.57
1:B:107:THR:O	1:B:108:ILE:HB	2.05	0.57
1:C:112:HIS:NE2	1:C:216:VAL:HG13	2.19	0.56
1:B:47:GLN:C	1:B:47:GLN:NE2	2.62	0.56
1:B:231:ALA:HB2	1:B:255:PHE:CD2	2.40	0.56
1:C:117:GLN:HB3	1:C:160:ARG:NH1	2.20	0.56
1:A:163:SER:HA	1:A:168:HIS:CD2	2.41	0.56
1:B:167:GLN:HE21	1:B:259:ARG:NH2	2.04	0.56
1:C:112:HIS:CE1	1:C:113:PRO:HD2	2.41	0.56
1:B:212:GLU:CG	1:B:220:THR:HB	2.36	0.55
1:C:59:ARG:O	1:C:62:GLU:HG2	2.06	0.55
1:A:139:TRP:HZ2	1:A:178:THR:HB	1.71	0.55
1:A:191:LEU:HD22	1:A:253:LEU:HD13	1.89	0.55
1:A:265:GLU:O	1:A:266:GLY:O	2.25	0.55
1:C:184:LEU:HA	1:C:187:ARG:CD	2.36	0.55
1:A:265:GLU:O	1:A:266:GLY:C	2.49	0.54
1:C:106:MET:O	1:C:107:THR:OG1	2.22	0.54
1:A:216:VAL:O	1:A:239:PHE:CE2	2.60	0.54
1:B:21:ARG:HH21	1:B:54:MET:HE3	1.72	0.54
1:B:183:PRO:HG2	1:B:186:GLU:HB2	1.88	0.54
1:C:117:GLN:NE2	1:C:160:ARG:HH12	2.04	0.54
1:A:15:THR:HG22	1:A:18:GLN:CG	2.38	0.54
1:B:108:ILE:HG13	1:B:112:HIS:H	1.73	0.54
1:B:277:LEU:HD22	2:B:473:HOH:O	2.07	0.54
1:C:19:TYR:CE1	1:C:32:LEU:HD22	2.43	0.54
1:B:21:ARG:HH21	1:B:54:MET:CE	2.20	0.54
1:C:140:ARG:HB2	2:C:392:HOH:O	2.07	0.54
1:A:58:LYS:HD2	1:A:288:ASN:CB	2.38	0.54
1:B:70:ARG:HH22	1:B:78:GLU:CD	2.16	0.54
1:C:231:ALA:HB2	1:C:255:PHE:CD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:THR:HA	1:B:128:ARG:HD3	1.90	0.53
1:C:220:THR:HG23	1:C:233:GLY:O	2.09	0.53
1:B:108:ILE:HD13	1:B:115:PHE:HD1	1.73	0.53
1:B:108:ILE:CG1	1:B:112:HIS:HB3	2.38	0.53
1:C:108:ILE:HD11	1:C:114:LEU:CB	2.30	0.53
1:C:190:ARG:O	1:C:194:MET:HG3	2.08	0.53
1:B:185:GLU:CD	1:B:185:GLU:H	2.10	0.53
1:C:122:ASP:OD1	1:C:125:LYS:HE2	2.08	0.53
1:A:21:ARG:HH21	1:A:54:MET:HE2	1.74	0.53
1:B:20:GLN:HA	1:B:20:GLN:NE2	2.24	0.53
1:B:108:ILE:HD12	1:B:114:LEU:CB	2.37	0.53
1:A:15:THR:HG22	1:A:18:GLN:CD	2.34	0.53
1:C:58:LYS:HE3	1:C:288:ASN:CB	2.37	0.52
1:A:19:TYR:HD1	1:A:32:LEU:HD21	1.74	0.52
1:C:120:TRP:CE3	1:C:126:CYS:HB3	2.45	0.52
1:B:107:THR:HB	1:B:240:LEU:HD21	1.90	0.52
1:A:42:ARG:HG2	2:A:315:HOH:O	2.10	0.52
1:B:217:TYR:HD1	1:B:217:TYR:O	1.93	0.52
1:B:236:GLY:HA2	1:B:237:PRO:C	2.34	0.52
1:A:19:TYR:CD1	1:A:32:LEU:HD21	2.45	0.51
1:A:267:THR:HG22	1:A:268:GLN:N	2.25	0.51
1:B:114:LEU:HD11	1:B:130:MET:HE1	1.91	0.51
1:B:212:GLU:HG2	1:B:220:THR:O	2.10	0.51
1:C:269:HIS:HB3	1:C:271:GLN:HG2	1.91	0.51
1:B:262:MET:HG3	1:B:267:THR:CG2	2.40	0.51
1:C:118:VAL:HG11	1:C:126:CYS:CB	2.41	0.51
1:B:31:GLN:O	1:B:34:MET:HG2	2.11	0.51
1:C:108:ILE:O	1:C:108:ILE:CG2	2.59	0.51
1:B:206:GLU:HG3	1:B:225:LYS:HD2	1.93	0.51
1:A:20:GLN:O	1:A:24:GLU:HG3	2.11	0.51
1:B:13:PHE:O	1:B:42:ARG:HD2	2.11	0.51
1:A:263:ILE:O	1:A:263:ILE:CG2	2.59	0.51
1:C:94:THR:HB	1:C:95:PRO:CD	2.41	0.51
1:B:105:LYS:C	1:B:107:THR:H	2.19	0.50
1:C:220:THR:HG22	1:C:221:VAL:N	2.26	0.50
1:A:29:GLY:O	1:A:33:GLU:HG3	2.11	0.50
1:C:129:PRO:CG	1:C:130:MET:HE2	2.32	0.50
1:A:70:ARG:HD3	2:A:358:HOH:O	2.12	0.50
1:B:167:GLN:OE1	1:B:262:MET:HE1	2.12	0.50
1:A:40:LEU:O	1:A:44:ARG:HG3	2.12	0.50
1:C:61:LEU:HD22	1:C:283:VAL:HG21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:SER:OG	1:C:168:HIS:HD2	1.95	0.50
1:B:40:LEU:C	1:B:42:ARG:N	2.66	0.49
1:C:15:THR:H	1:C:18:GLN:HE21	1.60	0.49
1:B:119:PHE:CZ	1:B:161:LYS:HB3	2.47	0.49
1:B:54:MET:HE2	1:B:287:ILE:CA	2.41	0.49
1:B:65:ARG:NH1	1:B:66:THR:HG23	2.27	0.49
1:C:108:ILE:HD13	1:C:115:PHE:CD2	2.48	0.49
1:A:94:THR:HB	1:A:95:PRO:CD	2.42	0.49
1:C:259:ARG:HG2	2:C:353:HOH:O	2.12	0.49
1:B:212:GLU:O	1:B:212:GLU:HG3	2.12	0.49
1:A:95:PRO:CD	1:B:277:LEU:HD13	2.42	0.49
1:B:143:GLU:HG2	1:B:148:LYS:HE3	1.94	0.49
1:C:59:ARG:HD3	1:C:60:HIS:N	2.27	0.49
1:C:160:ARG:NH1	1:C:162:GLU:O	2.46	0.49
1:B:54:MET:HE2	1:B:287:ILE:CB	2.42	0.49
1:B:108:ILE:HD13	1:B:115:PHE:CD1	2.48	0.49
1:B:206:GLU:O	1:B:225:LYS:HD3	2.13	0.49
1:A:215:VAL:HG13	1:A:216:VAL:N	2.26	0.49
1:B:114:LEU:HD23	1:B:114:LEU:O	2.13	0.49
1:B:228:LEU:CD1	1:B:263:ILE:HD12	2.38	0.49
1:B:284:ARG:CZ	1:B:286:ASN:HB2	2.43	0.48
1:B:20:GLN:HA	1:B:20:GLN:HE21	1.77	0.48
1:B:40:LEU:HD11	1:B:44:ARG:HE	1.78	0.48
1:B:262:MET:HA	1:B:267:THR:HG22	1.93	0.48
1:C:40:LEU:C	1:C:42:ARG:N	2.71	0.48
1:C:112:HIS:CE1	1:C:216:VAL:HG13	2.48	0.48
1:B:40:LEU:HD11	2:B:404:HOH:O	2.12	0.48
1:B:189:GLN:O	1:B:192:GLU:HB2	2.13	0.48
1:C:54:MET:CE	1:C:288:ASN:H	2.26	0.48
1:C:61:LEU:HD21	1:C:283:VAL:CG1	2.43	0.48
1:B:79:LYS:HD2	2:B:416:HOH:O	2.14	0.48
1:B:206:GLU:HG3	1:B:225:LYS:CD	2.44	0.48
1:A:15:THR:HG22	1:A:18:GLN:HB2	1.94	0.48
1:A:82:LYS:HE2	2:A:313:HOH:O	2.14	0.48
1:A:116:SER:O	1:A:161:LYS:HE2	2.15	0.47
1:A:106:MET:CE	1:A:244:TRP:NE1	2.78	0.47
1:B:21:ARG:NH2	1:B:54:MET:HE3	2.29	0.47
1:C:108:ILE:HA	1:C:112:HIS:HB2	1.97	0.47
1:C:118:VAL:HG11	1:C:126:CYS:SG	2.54	0.47
1:A:15:THR:HG22	1:A:18:GLN:NE2	2.30	0.47
1:A:216:VAL:O	1:A:239:PHE:HE2	1.98	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:GLN:NE2	1:B:259:ARG:NH2	2.63	0.47
1:B:215:VAL:HG22	1:B:215:VAL:O	2.15	0.47
1:B:65:ARG:HH11	1:B:65:ARG:HB3	1.79	0.46
1:C:168:HIS:HE1	1:C:170:ASN:OD1	1.97	0.46
1:A:51:HIS:HD2	2:A:335:HOH:O	1.97	0.46
1:A:209:LEU:HD12	1:A:223:VAL:HG22	1.97	0.46
1:A:269:HIS:CG	1:A:271:GLN:HG2	2.50	0.46
1:A:217:TYR:HB3	1:A:220:THR:OG1	2.15	0.46
1:A:272:SER:O	1:A:282:GLY:HA2	2.15	0.46
1:B:50:GLU:O	1:B:54:MET:HG3	2.16	0.46
1:B:209:LEU:HD23	1:B:223:VAL:HG22	1.96	0.46
1:A:33:GLU:HG2	2:A:374:HOH:O	2.16	0.46
1:A:229:GLU:HG3	1:A:229:GLU:O	2.14	0.46
1:B:108:ILE:HG23	1:B:115:PHE:CE1	2.51	0.46
1:B:228:LEU:HD13	1:B:263:ILE:CD1	2.38	0.46
1:C:68:LYS:O	1:C:69:HIS:HB2	2.14	0.46
1:B:20:GLN:O	1:B:24:GLU:HG3	2.16	0.46
1:A:118:VAL:CG1	1:A:126:CYS:HB2	2.46	0.46
1:A:122:ASP:HB3	1:A:124:LYS:H	1.82	0.45
1:A:146:TRP:CZ2	1:B:65:ARG:HB2	2.51	0.45
1:A:267:THR:CG2	1:A:268:GLN:N	2.79	0.45
1:B:108:ILE:HG13	1:B:112:HIS:CB	2.46	0.45
1:B:167:GLN:HE21	1:B:259:ARG:HH21	1.63	0.45
1:C:109:GLY:C	1:C:111:ASP:H	2.24	0.45
1:B:275:ARG:HG3	2:B:433:HOH:O	2.16	0.45
1:C:122:ASP:OD2	1:C:124:LYS:N	2.47	0.45
1:B:94:THR:HB	1:B:95:PRO:HD2	1.97	0.45
1:B:216:VAL:HG13	1:B:217:TYR:CD2	2.51	0.45
1:A:163:SER:CB	1:A:168:HIS:HB3	2.45	0.45
1:A:111:ASP:O	1:A:113:PRO:HD3	2.17	0.45
1:B:106:MET:HE1	1:B:136:TYR:CD2	2.51	0.45
1:B:116:SER:HA	2:B:434:HOH:O	2.16	0.45
1:C:159:TYR:CE2	1:C:171:GLU:HG2	2.52	0.45
1:C:167:GLN:HE22	1:C:269:HIS:HA	1.81	0.45
1:B:106:MET:HE2	1:B:244:TRP:NE1	2.30	0.45
1:C:107:THR:HG21	1:C:239:PHE:HE1	1.82	0.45
1:C:218:GLY:HA3	1:C:239:PHE:CE2	2.50	0.45
1:A:93:VAL:HG13	1:A:93:VAL:O	2.17	0.45
1:B:54:MET:HE2	1:B:287:ILE:CG2	2.44	0.45
1:B:182:THR:O	1:B:249:PRO:HB2	2.18	0.44
1:C:54:MET:HE2	1:C:288:ASN:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:LEU:O	1:C:80:LEU:HG	2.17	0.44
1:A:12:SER:HA	2:A:339:HOH:O	2.17	0.44
1:A:96:THR:HA	1:A:128:ARG:HD3	1.99	0.44
1:A:106:MET:CE	1:A:244:TRP:HE1	2.29	0.44
1:C:40:LEU:N	1:C:40:LEU:HD12	2.33	0.44
1:C:54:MET:HE2	1:C:288:ASN:H	1.81	0.44
1:C:108:ILE:HA	2:C:401:HOH:O	2.17	0.44
1:C:272:SER:OG	1:C:282:GLY:HA2	2.17	0.44
1:C:108:ILE:HG23	2:C:442:HOH:O	2.17	0.44
1:A:43:ASP:HA	2:A:315:HOH:O	2.17	0.44
1:A:190:ARG:HA	1:A:190:ARG:HH11	1.83	0.44
1:B:13:PHE:C	1:B:42:ARG:HD2	2.42	0.44
1:C:21:ARG:NH2	1:C:54:MET:CE	2.74	0.44
1:C:122:ASP:HA	2:C:360:HOH:O	2.17	0.44
1:A:120:TRP:CE3	1:A:126:CYS:HB3	2.52	0.44
1:B:122:ASP:OD2	1:B:122:ASP:C	2.61	0.44
1:A:40:LEU:HG	1:A:44:ARG:CZ	2.48	0.43
1:A:196:ARG:HD3	2:A:388:HOH:O	2.18	0.43
1:C:113:PRO:HD3	2:C:396:HOH:O	2.17	0.43
1:B:33:GLU:O	1:B:34:MET:C	2.62	0.43
1:B:108:ILE:CD1	1:B:112:HIS:HB3	2.49	0.43
1:C:185:GLU:OE1	1:C:185:GLU:N	2.28	0.43
1:C:73:LEU:HA	1:C:261:LEU:HD11	2.00	0.43
1:C:243:LYS:HB2	2:C:438:HOH:O	2.17	0.43
1:A:144:ARG:NH2	1:B:287:ILE:O	2.51	0.43
1:B:40:LEU:O	1:B:42:ARG:N	2.51	0.43
1:B:155:ILE:HG12	1:B:175:LEU:HD13	2.00	0.43
1:B:100:LYS:HE2	1:B:115:PHE:HE2	1.83	0.43
1:C:139:TRP:HZ2	1:C:178:THR:HB	1.84	0.43
1:C:192:GLU:OE2	1:C:209:LEU:HD13	2.18	0.43
1:A:159:TYR:CE2	1:A:171:GLU:HG2	2.54	0.43
1:A:216:VAL:O	1:A:218:GLY:N	2.52	0.43
1:C:54:MET:CE	1:C:287:ILE:HA	2.28	0.43
1:A:208:GLU:HG2	1:A:209:LEU:N	2.34	0.42
1:B:15:THR:C	1:B:17:VAL:H	2.26	0.42
1:B:62:GLU:OE1	1:B:65:ARG:NH1	2.52	0.42
1:B:118:VAL:HG13	1:B:126:CYS:HB2	2.01	0.42
1:B:155:ILE:HA	1:B:174:MET:O	2.18	0.42
1:C:122:ASP:OD2	1:C:122:ASP:C	2.62	0.42
1:A:161:LYS:HE3	2:A:385:HOH:O	2.18	0.42
1:B:103:LEU:HD12	1:B:115:PHE:HE1	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:106:MET:HE1	1:C:136:TYR:CD2	2.54	0.42
1:C:199:LEU:HD11	1:C:223:VAL:HG11	2.01	0.42
1:B:206:GLU:HG3	1:B:206:GLU:O	2.18	0.42
1:B:209:LEU:HD21	1:B:223:VAL:HG22	1.98	0.42
1:B:114:LEU:HD21	1:B:129:PRO:HG3	2.00	0.42
1:B:272:SER:HA	1:B:279:TYR:CG	2.55	0.42
1:C:271:GLN:NE2	2:C:463:HOH:O	2.52	0.42
1:A:60:HIS:HE1	1:A:281:ASP:OD2	2.02	0.42
1:A:19:TYR:HE1	1:A:32:LEU:HD11	1.82	0.42
1:B:217:TYR:O	1:B:217:TYR:CD1	2.71	0.42
1:B:117:GLN:NE2	1:B:162:GLU:OE2	2.53	0.42
1:B:231:ALA:HB2	1:B:255:PHE:CE2	2.54	0.42
1:A:90:VAL:O	1:A:152:ILE:HA	2.20	0.41
1:A:148:LYS:HA	1:A:149:PRO:C	2.45	0.41
1:B:20:GLN:HE21	1:B:20:GLN:CA	2.29	0.41
1:C:79:LYS:HE3	2:C:467:HOH:O	2.20	0.41
1:A:286:ASN:O	1:A:286:ASN:CG	2.63	0.41
1:B:40:LEU:O	1:B:41:SER:C	2.62	0.41
1:C:284:ARG:HG3	1:C:286:ASN:OD1	2.19	0.41
1:A:105:LYS:C	1:A:107:THR:H	2.29	0.41
1:A:191:LEU:HD23	1:A:191:LEU:HA	1.94	0.41
1:A:264:ARG:C	1:A:266:GLY:H	2.28	0.41
1:B:159:TYR:CD1	1:B:159:TYR:N	2.88	0.41
1:C:220:THR:CG2	1:C:221:VAL:N	2.83	0.41
1:B:40:LEU:C	1:B:42:ARG:H	2.29	0.41
1:B:106:MET:HE1	1:B:136:TYR:HD2	1.84	0.41
1:B:133:PRO:HG2	2:B:511:HOH:O	2.20	0.41
1:B:196:ARG:O	1:B:200:GLU:HB2	2.21	0.41
1:B:212:GLU:CD	1:B:220:THR:HB	2.46	0.41
1:A:105:LYS:HB2	1:A:244:TRP:CZ2	2.56	0.41
1:A:44:ARG:HG2	1:A:44:ARG:HH11	1.86	0.41
1:A:184:LEU:HG	2:A:318:HOH:O	2.20	0.41
1:B:108:ILE:HD11	1:B:115:PHE:H	1.86	0.41
1:A:76:LEU:O	1:A:80:LEU:HG	2.21	0.41
1:B:61:LEU:HD13	1:B:283:VAL:HG21	2.03	0.41
1:B:107:THR:O	1:B:107:THR:CG2	2.63	0.41
1:B:253:LEU:HD12	1:B:253:LEU:HA	1.92	0.41
1:C:40:LEU:C	1:C:42:ARG:H	2.28	0.41
1:C:16:LYS:HE2	1:C:20:GLN:OE1	2.21	0.41
1:B:206:GLU:CG	1:B:225:LYS:HE2	2.51	0.40
1:C:19:TYR:CD1	1:C:32:LEU:HD22	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:ARG:O	1:A:70:ARG:HG2	2.19	0.40
1:B:103:LEU:HD12	1:B:115:PHE:CD1	2.56	0.40
1:B:92:VAL:HG21	1:B:138:LEU:HD12	2.03	0.40
1:A:111:ASP:C	1:A:113:PRO:HD3	2.46	0.40
1:A:211:THR:HG22	1:A:212:GLU:N	2.37	0.40
1:A:106:MET:HE2	1:A:244:TRP:NE1	2.36	0.40
1:A:188:HIS:CE1	1:A:235:MET:HE1	2.56	0.40
1:B:117:GLN:HG2	1:B:162:GLU:CD	2.46	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	276/308 (90%)	249 (90%)	21 (8%)	6 (2%)	5	9
1	B	277/308 (90%)	253 (91%)	19 (7%)	5 (2%)	6	12
1	C	276/308 (90%)	264 (96%)	9 (3%)	3 (1%)	11	22
All	All	829/924 (90%)	766 (92%)	49 (6%)	14 (2%)	7	13

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	GLY
1	B	108	ILE
1	B	107	THR
1	B	218	GLY
1	B	286	ASN
1	C	111	ASP
1	A	217	TYR
1	A	268	GLN
1	A	210	VAL

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Mol	Chain	Res	Type
1	A	16	LYS
1	B	111	ASP
1	C	107	THR
1	A	263	ILE
1	C	237	PRO

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	242/269 (90%)	227 (94%)	15 (6%)	16	34
1	B	243/269 (90%)	224 (92%)	19 (8%)	11	24
1	C	242/269 (90%)	230 (95%)	12 (5%)	22	44
All	All	727/807 (90%)	681 (94%)	46 (6%)	16	34

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	61	LEU
1	A	67	VAL
1	A	70	ARG
1	A	86	GLN
1	A	100	LYS
1	A	103	LEU
1	A	114	LEU
1	A	117	GLN
1	A	136	TYR
1	A	175	LEU
1	A	185	GLU
1	A	222	ASP
1	A	260	LEU
1	A	288	ASN
1	B	17	VAL
1	B	47	GLN

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Mol	Chain	Res	Type
1	B	61	LEU
1	B	65	ARG
1	B	67	VAL
1	B	70	ARG
1	B	103	LEU
1	B	107	THR
1	B	118	VAL
1	B	130	MET
1	B	136	TYR
1	B	140	ARG
1	B	175	LEU
1	B	185	GLU
1	B	242	GLU
1	B	253	LEU
1	B	260	LEU
1	B	277	LEU
1	B	285	LEU
1	C	22	LEU
1	C	47	GLN
1	C	59	ARG
1	C	61	LEU
1	C	65	ARG
1	C	70	ARG
1	C	103	LEU
1	C	108	ILE
1	C	136	TYR
1	C	144	ARG
1	C	185	GLU
1	C	253	LEU

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

Mol	Chain	Res	Type
1	A	18	GLN
1	A	52	GLN
1	A	56	GLN
1	A	60	HIS
1	A	63	GLN
1	A	117	GLN
1	A	164	GLN
1	A	188	HIS
1	A	189	GLN

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Mol	Chain	Res	Type
1	A	271	GLN
1	A	288	ASN
1	B	18	GLN
1	B	20	GLN
1	B	26	ASN
1	B	47	GLN
1	B	52	GLN
1	B	167	GLN
1	B	176	ASN
1	B	271	GLN
1	B	286	ASN
1	B	288	ASN
1	C	18	GLN
1	C	26	ASN
1	C	47	GLN
1	C	52	GLN
1	C	56	GLN
1	C	112	HIS
1	C	117	GLN
1	C	167	GLN
1	C	168	HIS
1	C	176	ASN
1	C	271	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 ⓘ

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	278/308 (90%)	0.42	37 (13%) 7 6	17, 37, 77, 86	0
1	B	279/308 (90%)	0.31	27 (9%) 13 11	14, 36, 78, 88	0
1	C	278/308 (90%)	0.09	12 (4%) 40 35	11, 34, 66, 78	0
All	All	835/924 (90%)	0.27	76 (9%) 15 13	11, 36, 76, 88	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215	VAL	8.3
1	C	109	GLY	7.5
1	B	216	VAL	6.9
1	C	113	PRO	5.2
1	B	217	TYR	5.1
1	A	216	VAL	5.0
1	B	218	GLY	4.9
1	B	165	GLY	4.9
1	C	112	HIS	4.0
1	A	19	TYR	3.9
1	A	166	ALA	3.8
1	B	219	ASP	3.6
1	A	217	TYR	3.5
1	B	214	SER	3.5
1	A	263	ILE	3.5
1	C	108	ILE	3.4
1	C	111	ASP	3.3
1	B	10	LEU	3.3
1	A	220	THR	3.2
1	A	213	SER	3.2
1	B	213	SER	3.2
1	B	162	GLU	3.2
1	A	32	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	211	THR	3.2
1	C	36	PHE	3.1
1	C	62	GLU	3.1
1	B	163	SER	3.1
1	B	108	ILE	3.1
1	B	220	THR	3.0
1	B	112	HIS	3.0
1	A	269	HIS	3.0
1	A	13	PHE	3.0
1	B	109	GLY	2.9
1	A	265	GLU	2.9
1	A	163	SER	2.9
1	A	165	GLY	2.9
1	A	164	GLN	2.9
1	A	227	ASP	2.8
1	A	29	GLY	2.8
1	B	40	LEU	2.8
1	C	35	GLY	2.7
1	A	268	GLN	2.7
1	B	47	GLN	2.7
1	B	107	THR	2.7
1	A	288	ASN	2.6
1	A	108	ILE	2.6
1	B	44	ARG	2.6
1	B	111	ASP	2.5
1	B	39	ALA	2.5
1	A	167	GLN	2.5
1	A	36	PHE	2.5
1	B	164	GLN	2.4
1	B	113	PRO	2.4
1	A	162	GLU	2.4
1	A	215	VAL	2.4
1	B	37	SER	2.4
1	B	166	ALA	2.4
1	A	11	SER	2.3
1	A	30	GLU	2.3
1	A	168	HIS	2.3
1	C	239	PHE	2.3
1	A	39	ALA	2.2
1	A	107	THR	2.2
1	A	274	ALA	2.2
1	C	34	MET	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	228	LEU	2.2
1	A	273	MET	2.1
1	A	109	GLY	2.1
1	A	34	MET	2.1
1	C	110	GLU	2.1
1	A	209	LEU	2.0
1	B	110	GLU	2.0
1	C	288	ASN	2.0
1	A	219	ASP	2.0
1	A	232	SER	2.0
1	B	11	SER	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.