



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:52 PM UTC

PDB ID : 3ESH / pdb_00003esh
Title : Crystal structure of a probable metal-dependent hydrolase from *Staphylococcus aureus*. Northeast Structural Genomics target ZR314
Authors : Seetharaman, J.; Chen, Y.; Wang, H.; Janjua, H.; Foote, E.L.; Xiao, R.; Nair, R.; Everett, J.K.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2008-10-06
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
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

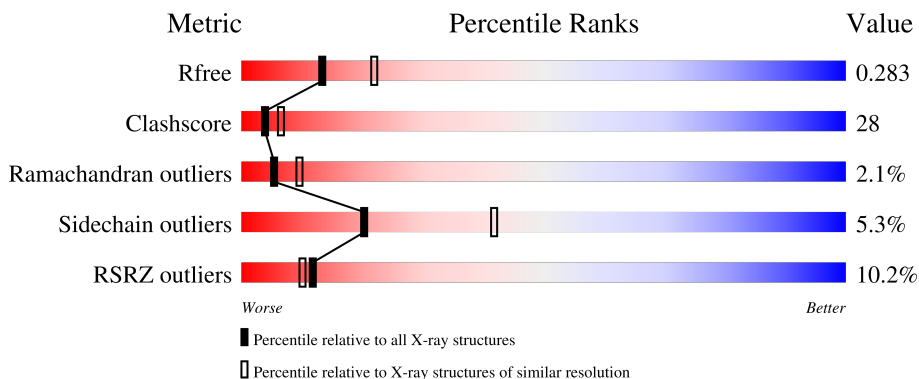
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

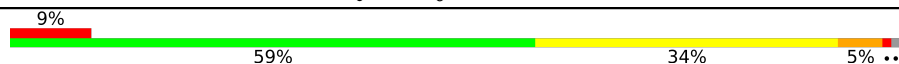


The reported resolution of this entry is 2.50 Å.

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	280	
1	B	280	
1	C	280	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6886 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 Protein similar to metal-dependent hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	277	2250	1441	381	419	9	0	0	0
1	B	275	2235	1430	379	417	9	0	0	0
1	C	275	2235	1430	379	417	9	0	0	0

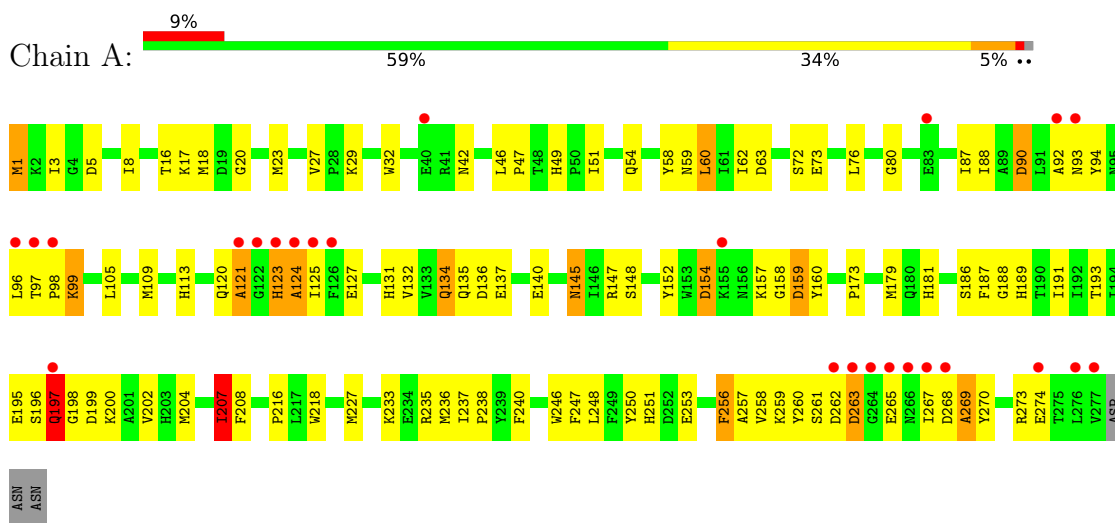
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	64	Total 64	O 64	0	0
2	B	42	Total 42	O 42	0	0
2	C	60	Total 60	O 60	0	0

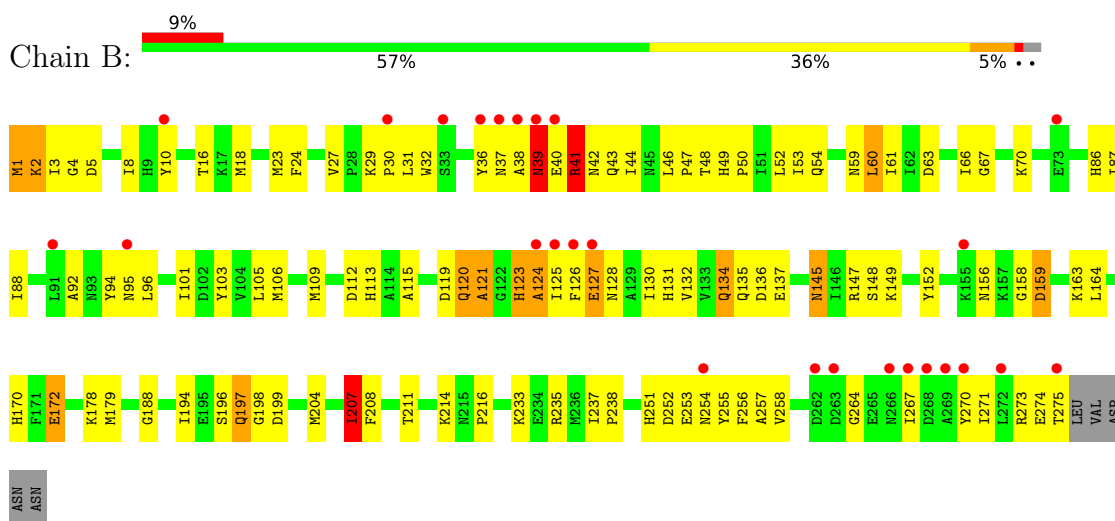
3 Residue-property plots [i](#)

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: Protein similar to metal-dependent hydrolase

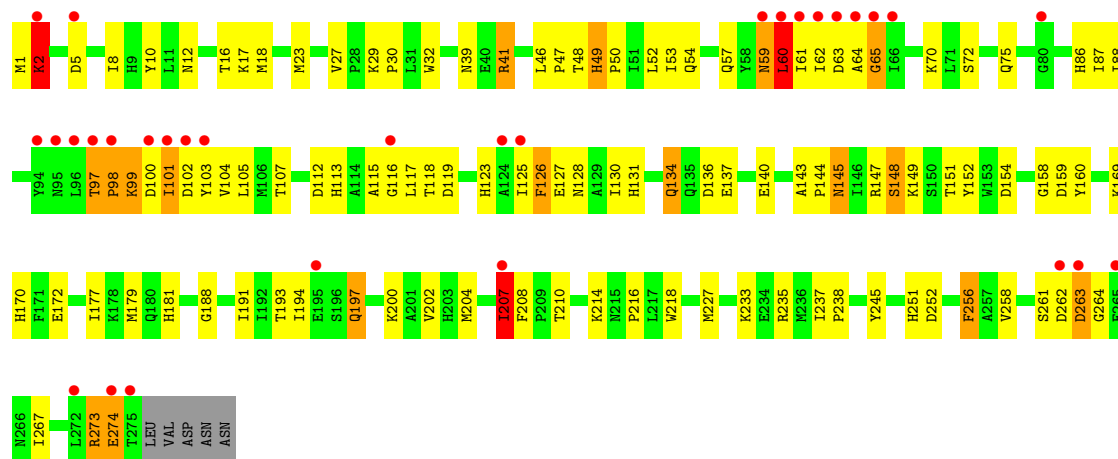


- Molecule 1: Protein similar to metal-dependent hydrolase



- Molecule 1: Protein similar to metal-dependent hydrolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	113.11Å 113.11Å 155.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.77 – 2.50 45.77 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.0 (45.77-2.50) 93.7 (45.77-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.45 (at 2.39Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.228 , 0.274 0.242 , 0.283	Depositor DCC
R_{free} test set	3098 reflections (4.13%)	wwPDB-VP
Wilson B-factor (Å ²)	39.9	Xtrriage
Anisotropy	0.478	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 33.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6886	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

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	0.53	0/2307	0.99	14/3116 (0.4%)
1	B	0.50	0/2292	1.00	15/3095 (0.5%)
1	C	0.52	0/2292	0.94	15/3095 (0.5%)
All	All	0.52	0/6891	0.98	44/9306 (0.5%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	125	ILE	N-CA-C	-11.46	97.26	110.21
1	A	125	ILE	N-CA-C	-11.35	97.38	110.21
1	B	41	ARG	N-CA-C	-8.64	102.74	113.28
1	A	127	GLU	N-CA-C	7.50	120.33	111.71
1	A	121	ALA	N-CA-C	-7.07	103.01	113.40
1	B	127	GLU	N-CA-C	-7.06	101.02	110.55
1	A	269	ALA	N-CA-C	7.03	119.57	108.67
1	C	126	PHE	N-CA-C	6.87	118.23	108.54
1	B	121	ALA	N-CA-C	-6.75	104.99	113.23
1	A	197	GLN	CB-CA-C	-6.73	108.83	116.63
1	A	158	GLY	N-CA-C	6.56	120.14	110.63
1	B	158	GLY	N-CA-C	6.41	119.97	110.42
1	C	256	PHE	N-CA-C	-6.36	104.47	112.93
1	A	109	MSE	N-CA-C	6.33	119.93	112.72
1	A	207	ILE	N-CA-C	-6.32	105.44	113.22
1	C	207	ILE	N-CA-C	-6.01	105.83	113.22
1	C	41	ARG	N-CA-C	-5.99	105.74	113.16
1	B	86	HIS	N-CA-C	-5.93	104.24	112.30
1	B	207	ILE	N-CA-C	-5.83	106.05	113.22
1	C	86	HIS	N-CA-C	-5.82	104.85	112.41
1	B	39	ASN	N-CA-C	-5.78	98.48	110.80
1	C	197	GLN	CB-CA-C	-5.67	110.06	116.63
1	B	135	GLN	N-CA-C	5.58	120.53	111.37
1	C	154	ASP	N-CA-C	-5.58	105.20	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	109	MSE	N-CA-C	5.50	118.81	112.97
1	C	2	LYS	N-CA-C	5.49	117.68	108.73
1	B	208	PHE	CA-C-N	5.42	126.61	119.84
1	B	208	PHE	C-N-CA	5.42	126.61	119.84
1	B	172	GLU	CA-C-N	5.35	124.81	119.24
1	B	172	GLU	C-N-CA	5.35	124.81	119.24
1	A	256	PHE	N-CA-C	-5.34	106.89	113.41
1	C	273	ARG	N-CA-C	5.33	117.77	108.76
1	C	208	PHE	CA-C-N	5.25	126.40	119.84
1	C	208	PHE	C-N-CA	5.25	126.40	119.84
1	C	172	GLU	CA-C-N	5.24	126.39	119.84
1	C	172	GLU	C-N-CA	5.24	126.39	119.84
1	A	208	PHE	CA-C-N	5.23	126.38	119.84
1	A	208	PHE	C-N-CA	5.23	126.38	119.84
1	A	154	ASP	N-CA-C	-5.22	105.59	111.28
1	A	135	GLN	N-CA-C	5.16	116.90	111.28
1	A	58	TYR	N-CA-C	5.15	118.12	110.14
1	C	49	HIS	N-CA-C	5.13	118.65	108.59
1	C	107	THR	N-CA-C	-5.12	105.34	112.45
1	B	120	GLN	N-CA-C	-5.02	104.23	110.41

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	2250	0	2149	133	0
1	B	2235	0	2129	107	0
1	C	2235	0	2129	138	0
2	A	64	0	0	0	0
2	B	42	0	0	2	0
2	C	60	0	0	1	0
All	All	6886	0	6407	372	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 (372) 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:258:VAL:HG22	1:A:267:ILE:CD1	1.69	1.22
1:C:97:THR:HG23	1:C:98:PRO:HD2	1.18	1.16
1:C:60:LEU:H	1:C:60:LEU:HD23	1.09	1.09
1:C:60:LEU:H	1:C:60:LEU:CD2	1.64	1.09
1:A:267:ILE:HG21	1:A:270:TYR:OH	1.53	1.05
1:C:60:LEU:HD23	1:C:60:LEU:N	1.70	1.05
1:C:52:LEU:CD1	1:C:61:ILE:HD11	1.86	1.05
1:C:97:THR:HG23	1:C:98:PRO:CD	1.90	1.01
1:A:258:VAL:CG2	1:A:267:ILE:HD11	1.94	0.97
1:A:258:VAL:HG22	1:A:267:ILE:HD11	0.98	0.95
1:B:60:LEU:HD22	1:B:103:TYR:HB2	1.49	0.94
1:B:113:HIS:HE1	2:B:281:HOH:O	1.50	0.93
1:C:52:LEU:HD13	1:C:61:ILE:HD11	1.47	0.93
1:C:52:LEU:HA	1:C:61:ILE:CD1	1.99	0.91
1:A:258:VAL:HG21	1:A:267:ILE:HG12	1.50	0.91
1:C:63:ASP:OD2	1:C:113:HIS:HD2	1.53	0.91
1:C:97:THR:CG2	1:C:98:PRO:HD2	2.02	0.87
1:A:258:VAL:CG2	1:A:267:ILE:CD1	2.50	0.86
1:A:258:VAL:CG2	1:A:267:ILE:HG12	2.06	0.86
1:A:260:TYR:CD1	1:A:267:ILE:HB	2.11	0.85
1:A:16:THR:HB	1:A:18:MSE:HE3	1.56	0.85
1:A:258:VAL:CG2	1:A:267:ILE:CG1	2.53	0.85
1:C:134:GLN:HE21	1:C:136:ASP:HB2	1.41	0.85
1:B:145:ASN:HD22	1:B:147:ARG:H	1.25	0.84
1:C:145:ASN:HD22	1:C:147:ARG:H	1.23	0.84
1:C:2:LYS:H	1:C:2:LYS:HD3	1.43	0.83
1:C:54:GLN:HG2	1:C:59:ASN:HB3	1.60	0.83
1:A:120:GLN:O	1:A:121:ALA:HB3	1.79	0.83
1:A:134:GLN:NE2	1:A:136:ASP:H	1.76	0.82
1:B:16:THR:HB	1:B:18:MSE:HE3	1.61	0.82
1:C:123:HIS:HA	1:C:159:ASP:OD2	1.79	0.82
1:B:134:GLN:HE21	1:B:137:GLU:H	1.25	0.82
1:C:52:LEU:HD12	1:C:61:ILE:HD11	1.62	0.81
1:B:105:LEU:HD22	1:B:179:MSE:SE	2.30	0.81
1:B:134:GLN:NE2	1:B:136:ASP:N	2.29	0.81
1:C:60:LEU:HB3	1:C:103:TYR:HB2	1.64	0.80
1:A:227:MSE:HE1	1:C:147:ARG:HD3	1.64	0.79
1:B:172:GLU:HG3	1:B:178:LYS:HG2	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:GLN:NE2	1:B:96:LEU:HD11	1.97	0.79
1:A:260:TYR:CE1	1:A:267:ILE:HB	2.18	0.79
1:B:41:ARG:HB2	1:B:43:GLN:HG3	1.63	0.79
1:A:147:ARG:HD3	1:C:227:MSE:HE1	1.64	0.78
1:C:52:LEU:HA	1:C:61:ILE:HD12	1.64	0.77
1:B:134:GLN:NE2	1:B:136:ASP:H	1.83	0.76
1:B:48:THR:O	1:B:50:PRO:HD3	1.86	0.75
1:C:134:GLN:NE2	1:C:136:ASP:HB2	2.00	0.75
1:A:145:ASN:HD22	1:A:147:ARG:H	1.35	0.75
1:B:47:PRO:HB3	1:B:49:HIS:CE1	2.21	0.75
1:B:134:GLN:HE22	1:B:136:ASP:H	1.35	0.74
1:A:134:GLN:HE21	1:A:136:ASP:HB2	1.50	0.74
1:C:60:LEU:O	1:C:61:ILE:HD13	1.87	0.74
1:C:64:ALA:O	1:C:116:GLY:HA3	1.89	0.73
1:C:123:HIS:HE1	1:C:127:GLU:OE1	1.72	0.73
1:C:134:GLN:HG3	1:C:137:GLU:H	1.54	0.73
1:A:145:ASN:ND2	1:A:148:SER:H	1.87	0.72
1:B:38:ALA:C	1:B:39:ASN:O	2.25	0.72
1:A:123:HIS:O	1:A:159:ASP:HB2	1.90	0.71
1:A:267:ILE:CG2	1:A:270:TYR:OH	2.34	0.71
1:A:227:MSE:HE1	1:C:147:ARG:CD	2.20	0.71
1:B:214:LYS:NZ	1:B:264:GLY:O	2.22	0.71
1:A:147:ARG:CD	1:C:227:MSE:HE1	2.20	0.71
1:C:103:TYR:CE1	1:C:130:ILE:HD12	2.26	0.70
1:A:120:GLN:O	1:A:121:ALA:CB	2.40	0.70
1:A:29:LYS:HA	1:A:32:TRP:NE1	2.07	0.70
1:A:123:HIS:O	1:A:124:ALA:HB2	1.92	0.70
1:C:258:VAL:HG13	1:C:267:ILE:HG23	1.74	0.70
1:C:103:TYR:CD1	1:C:130:ILE:HB	2.26	0.69
1:C:214:LYS:NZ	1:C:264:GLY:O	2.24	0.69
1:A:134:GLN:NE2	1:A:136:ASP:HB2	2.06	0.69
1:B:127:GLU:O	1:B:128:ASN:HB2	1.91	0.68
1:A:147:ARG:NE	1:C:227:MSE:HE1	2.08	0.68
1:B:145:ASN:ND2	1:B:148:SER:H	1.91	0.68
1:A:258:VAL:HG22	1:A:267:ILE:CG1	2.21	0.68
1:C:102:ASP:O	1:C:103:TYR:CD2	2.46	0.68
1:C:63:ASP:OD2	1:C:113:HIS:CD2	2.42	0.68
1:A:261:SER:O	1:A:263:ASP:N	2.27	0.67
1:C:50:PRO:HG3	1:C:65:GLY:H	1.58	0.67
1:C:52:LEU:HA	1:C:61:ILE:HD13	1.76	0.67
1:A:267:ILE:HG21	1:A:270:TYR:HH	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ALA:O	1:B:39:ASN:O	2.13	0.66
1:B:41:ARG:CZ	1:B:41:ARG:HB3	2.24	0.66
1:A:16:THR:HB	1:A:18:MSE:CE	2.26	0.66
1:C:145:ASN:ND2	1:C:147:ARG:H	1.93	0.66
1:A:63:ASP:OD2	1:A:113:HIS:HD2	1.78	0.66
1:B:123:HIS:O	1:B:124:ALA:HB2	1.95	0.66
1:C:103:TYR:CE1	1:C:130:ILE:HB	2.31	0.65
1:A:29:LYS:HA	1:A:32:TRP:CD1	2.31	0.65
1:C:105:LEU:HB3	1:C:179:MSE:HE3	1.77	0.65
1:C:49:HIS:HB2	1:C:273:ARG:NH2	2.12	0.65
1:C:102:ASP:O	1:C:103:TYR:CG	2.50	0.65
1:B:23:MSE:HE2	1:B:46:LEU:HD11	1.77	0.64
1:C:17:LYS:C	1:C:18:MSE:HE2	2.23	0.64
1:C:97:THR:O	1:C:99:LYS:N	2.30	0.64
1:A:5:ASP:CG	1:A:197:GLN:HE22	2.05	0.64
1:A:97:THR:O	1:A:99:LYS:N	2.30	0.63
1:C:2:LYS:H	1:C:2:LYS:CD	2.10	0.63
1:C:134:GLN:NE2	1:C:136:ASP:H	1.97	0.63
1:C:145:ASN:ND2	1:C:148:SER:H	1.97	0.63
1:A:227:MSE:HE1	1:C:147:ARG:NE	2.14	0.63
1:C:52:LEU:HD12	1:C:61:ILE:CD1	2.29	0.62
1:B:5:ASP:CG	1:B:197:GLN:HE22	2.06	0.62
1:B:128:ASN:H	1:B:163:LYS:HZ2	1.45	0.62
1:C:10:TYR:OH	1:C:274:GLU:N	2.33	0.61
1:C:103:TYR:CZ	1:C:130:ILE:HD12	2.35	0.61
1:A:181:HIS:HE1	1:A:187:PHE:O	1.83	0.61
1:A:261:SER:HB3	1:A:268:ASP:OD1	2.00	0.61
1:B:123:HIS:O	1:B:159:ASP:HB2	2.00	0.61
1:B:29:LYS:HA	1:B:32:TRP:NE1	2.15	0.61
1:B:148:SER:HB3	1:B:152:TYR:CD2	2.36	0.61
1:A:1:MSE:HE3	1:A:8:ILE:HB	1.83	0.60
1:B:145:ASN:ND2	1:B:147:ARG:H	1.96	0.60
1:B:258:VAL:HG13	1:B:267:ILE:HG23	1.84	0.59
1:A:145:ASN:HD21	1:A:148:SER:H	1.51	0.59
1:C:148:SER:HB3	1:C:152:TYR:HD2	1.66	0.59
1:A:124:ALA:HB2	1:A:159:ASP:HB2	1.84	0.59
1:A:259:LYS:C	1:A:267:ILE:HD12	2.28	0.58
1:C:47:PRO:HD2	1:C:251:HIS:O	2.03	0.58
1:A:260:TYR:CD1	1:A:267:ILE:HD13	2.38	0.58
1:A:148:SER:HB2	1:A:152:TYR:CD2	2.39	0.58
1:A:260:TYR:HA	1:A:267:ILE:HD12	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:MSE:HE3	1:C:2:LYS:HE3	1.86	0.57
1:C:87:ILE:HG23	1:C:88:ILE:N	2.19	0.57
1:A:134:GLN:HE21	1:A:136:ASP:CB	2.16	0.57
1:B:120:GLN:O	1:B:121:ALA:HB3	2.04	0.57
1:C:134:GLN:HG2	1:C:188:GLY:O	2.04	0.57
1:B:70:LYS:NZ	1:B:156:ASN:ND2	2.52	0.56
1:A:1:MSE:HG3	1:A:256:PHE:HE2	1.70	0.56
1:B:3:ILE:HG23	1:B:199:ASP:CB	2.35	0.56
1:B:41:ARG:O	1:B:42:ASN:HB3	2.06	0.56
1:C:52:LEU:CD1	1:C:61:ILE:CD1	2.74	0.56
1:A:49:HIS:HD2	1:A:273:ARG:NE	2.04	0.56
1:C:140:GLU:HG2	1:C:152:TYR:CE2	2.41	0.55
1:B:254:ASN:O	1:B:273:ARG:N	2.34	0.55
1:B:101:ILE:HG13	1:B:126:PHE:CE2	2.41	0.55
1:C:193:THR:HG22	1:C:202:VAL:HG23	1.89	0.55
1:A:3:ILE:HG23	1:A:199:ASP:HB2	1.89	0.55
1:B:10:TYR:HB3	1:B:256:PHE:CE1	2.41	0.55
1:C:52:LEU:CA	1:C:61:ILE:CD1	2.82	0.55
1:C:97:THR:CG2	1:C:98:PRO:CD	2.72	0.55
1:B:134:GLN:HE21	1:B:137:GLU:N	2.01	0.54
1:B:59:ASN:C	1:B:60:LEU:HD23	2.32	0.54
1:B:41:ARG:O	1:B:42:ASN:CB	2.55	0.54
1:C:181:HIS:HE1	1:C:188:GLY:HA2	1.72	0.54
1:A:62:ILE:O	1:A:105:LEU:O	2.26	0.54
1:A:259:LYS:C	1:A:267:ILE:CD1	2.81	0.54
1:A:267:ILE:HG12	1:A:270:TYR:CZ	2.42	0.54
1:C:134:GLN:HE21	1:C:136:ASP:CB	2.16	0.54
1:B:94:TYR:O	1:B:95:ASN:CG	2.50	0.54
1:A:134:GLN:HG3	1:A:137:GLU:H	1.73	0.54
1:B:4:GLY:HA3	1:B:196:SER:HB3	1.90	0.54
1:B:134:GLN:HE22	1:B:136:ASP:N	1.98	0.54
1:B:128:ASN:N	1:B:163:LYS:HZ2	2.06	0.54
1:A:47:PRO:HD2	1:A:251:HIS:O	2.08	0.54
1:C:99:LYS:HG3	1:C:128:ASN:ND2	2.22	0.54
1:B:131:HIS:O	1:B:164:LEU:HD12	2.08	0.53
1:C:2:LYS:HD3	1:C:2:LYS:N	2.17	0.53
1:C:39:ASN:OD1	1:C:41:ARG:N	2.41	0.53
1:A:27:VAL:HG21	1:A:218:TRP:CD2	2.44	0.53
1:A:134:GLN:NE2	1:A:136:ASP:CB	2.71	0.53
1:A:123:HIS:O	1:A:124:ALA:CB	2.56	0.53
1:B:29:LYS:HA	1:B:32:TRP:CD1	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:LEU:CD2	1:C:60:LEU:N	2.38	0.53
1:A:23:MSE:HE2	1:A:46:LEU:HD11	1.91	0.53
1:A:181:HIS:HE1	1:A:188:GLY:HA2	1.74	0.53
1:C:1:MSE:N	1:C:8:ILE:O	2.41	0.53
1:C:16:THR:HB	1:C:18:MSE:HE3	1.90	0.53
1:C:53:ILE:O	1:C:59:ASN:HA	2.09	0.53
1:A:1:MSE:CE	1:A:248:LEU:HD21	2.39	0.52
1:A:134:GLN:HG2	1:A:188:GLY:O	2.08	0.52
1:A:140:GLU:HG2	1:A:152:TYR:CE2	2.44	0.52
1:C:48:THR:O	1:C:50:PRO:HD3	2.09	0.52
1:C:131:HIS:CE1	1:C:160:TYR:HB2	2.44	0.52
1:A:260:TYR:CG	1:A:267:ILE:HD13	2.44	0.52
1:A:181:HIS:CE1	1:A:188:GLY:HA2	2.45	0.52
1:C:101:ILE:CG1	1:C:126:PHE:CD2	2.93	0.52
1:C:145:ASN:HD22	1:C:145:ASN:C	2.18	0.52
1:A:186:SER:HB3	1:A:189:HIS:HB2	1.92	0.52
1:A:145:ASN:ND2	1:A:147:ARG:H	2.05	0.51
1:B:123:HIS:O	1:B:124:ALA:CB	2.58	0.51
1:B:46:LEU:HD22	1:B:251:HIS:HB3	1.91	0.51
1:C:52:LEU:CA	1:C:61:ILE:HD12	2.39	0.51
1:B:47:PRO:HG2	1:B:253:GLU:HA	1.93	0.51
1:C:53:ILE:HD11	1:C:194:ILE:HD11	1.93	0.51
1:A:60:LEU:N	1:A:60:LEU:HD23	2.24	0.51
1:A:145:ASN:HD22	1:A:145:ASN:C	2.19	0.51
1:B:16:THR:HB	1:B:18:MSE:CE	2.38	0.50
1:B:47:PRO:HD2	1:B:251:HIS:O	2.11	0.50
1:B:63:ASP:OD2	1:B:113:HIS:HD2	1.94	0.50
1:B:3:ILE:HG23	1:B:199:ASP:HB3	1.92	0.50
1:A:88:ILE:HD12	1:A:88:ILE:N	2.27	0.50
1:A:191:ILE:HD13	1:A:236:MSE:CE	2.41	0.50
1:C:127:GLU:O	1:C:128:ASN:HB2	2.12	0.50
1:C:47:PRO:O	1:C:251:HIS:HA	2.11	0.50
1:A:258:VAL:HG21	1:A:267:ILE:CG1	2.26	0.50
1:B:92:ALA:C	1:B:94:TYR:H	2.18	0.50
1:A:247:PHE:O	1:A:257:ALA:HA	2.12	0.50
1:C:134:GLN:NE2	1:C:136:ASP:CB	2.74	0.50
1:A:237:ILE:HB	1:A:238:PRO:HD3	1.94	0.49
1:C:10:TYR:HB3	1:C:256:PHE:CE1	2.47	0.49
1:C:64:ALA:O	1:C:65:GLY:C	2.55	0.49
1:B:61:ILE:CG1	1:B:101:ILE:HD12	2.42	0.49
1:A:145:ASN:HD22	1:A:147:ARG:N	2.09	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:HIS:CE1	1:A:187:PHE:O	2.64	0.49
1:A:259:LYS:O	1:A:267:ILE:HD12	2.12	0.49
1:A:267:ILE:HG23	1:A:270:TYR:HE2	1.77	0.49
1:B:29:LYS:HB3	1:B:30:PRO:HD3	1.93	0.49
1:B:124:ALA:HB2	1:B:159:ASP:HB2	1.94	0.49
1:C:23:MSE:HE2	1:C:46:LEU:HD11	1.95	0.49
1:A:267:ILE:HG23	1:A:270:TYR:CE2	2.48	0.49
1:B:49:HIS:HB2	1:B:273:ARG:HH21	1.78	0.49
1:C:125:ILE:HG22	1:C:125:ILE:O	2.13	0.49
1:B:128:ASN:HA	1:B:163:LYS:HZ2	1.77	0.49
1:B:132:VAL:HG11	1:B:179:MSE:SE	2.63	0.49
1:B:119:ASP:OD1	1:B:120:GLN:O	2.30	0.49
1:C:100:ASP:O	1:C:101:ILE:C	2.56	0.49
1:A:16:THR:CB	1:A:18:MSE:HE3	2.36	0.48
1:A:191:ILE:CD1	1:A:236:MSE:HE1	2.43	0.48
1:A:97:THR:HG23	1:A:98:PRO:HD2	1.94	0.48
1:C:104:VAL:HG11	1:C:117:LEU:HD21	1.96	0.48
1:A:267:ILE:CG1	1:A:270:TYR:CE2	2.96	0.48
1:B:29:LYS:C	1:B:31:LEU:H	2.22	0.48
1:C:181:HIS:CE1	1:C:188:GLY:HA2	2.48	0.48
1:B:128:ASN:HA	1:B:163:LYS:NZ	2.29	0.48
1:C:29:LYS:HB3	1:C:30:PRO:HD3	1.96	0.48
1:A:134:GLN:NE2	1:A:136:ASP:N	2.53	0.48
1:C:12:ASN:HA	1:C:49:HIS:ND1	2.29	0.48
1:B:61:ILE:HG12	1:B:101:ILE:HD12	1.95	0.48
1:C:60:LEU:H	1:C:60:LEU:HD22	1.70	0.48
1:A:3:ILE:HG23	1:A:199:ASP:CB	2.44	0.47
1:B:1:MSE:HE2	1:B:8:ILE:HB	1.96	0.47
1:B:36:TYR:HD2	1:B:44:ILE:HB	1.79	0.47
1:C:143:ALA:N	1:C:144:PRO:HD3	2.29	0.47
1:A:29:LYS:HG3	1:A:32:TRP:CZ2	2.49	0.47
1:B:211:THR:HG21	1:B:255:TYR:CD2	2.49	0.47
1:B:216:PRO:HA	1:B:233:LYS:HE2	1.97	0.47
1:C:112:ASP:OD2	1:C:113:HIS:CE1	2.67	0.47
1:B:237:ILE:HB	1:B:238:PRO:HD3	1.96	0.47
1:C:70:LYS:HG3	1:C:115:ALA:O	2.14	0.47
1:A:46:LEU:HD22	1:A:251:HIS:HB3	1.96	0.47
1:B:70:LYS:HZ3	1:B:156:ASN:ND2	2.11	0.47
1:B:92:ALA:C	1:B:94:TYR:N	2.72	0.47
1:C:145:ASN:H	1:C:148:SER:HB2	1.79	0.47
1:C:237:ILE:HB	1:C:238:PRO:HD3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:LEU:HB3	1:A:179:MSE:HE3	1.97	0.47
1:C:99:LYS:HG3	1:C:128:ASN:HD22	1.80	0.47
1:A:148:SER:HB2	1:A:152:TYR:CE2	2.50	0.46
1:B:257:ALA:HB3	1:B:271:ILE:HD12	1.97	0.46
1:C:29:LYS:HA	1:C:32:TRP:NE1	2.30	0.46
1:B:70:LYS:HZ1	1:B:156:ASN:HD22	1.62	0.46
1:B:145:ASN:HD22	1:B:147:ARG:N	2.04	0.46
1:C:57:GLN:HG2	2:C:339:HOH:O	2.14	0.46
1:A:154:ASP:OD2	1:A:157:LYS:NZ	2.49	0.46
1:A:204:MSE:HB3	1:A:207:ILE:HG22	1.98	0.46
1:A:191:ILE:CD1	1:A:236:MSE:CE	2.93	0.46
1:A:47:PRO:HG2	1:A:253:GLU:HA	1.96	0.46
1:A:132:VAL:HG11	1:A:179:MSE:SE	2.65	0.46
1:B:145:ASN:HD22	1:B:145:ASN:C	2.22	0.46
1:A:17:LYS:C	1:A:18:MSE:HE2	2.41	0.46
1:C:252:ASP:O	1:C:273:ARG:CD	2.64	0.46
1:B:53:ILE:HD13	1:B:194:ILE:HG12	1.98	0.46
1:B:128:ASN:CA	1:B:163:LYS:HZ2	2.28	0.46
1:B:134:GLN:NE2	1:B:136:ASP:HB2	2.30	0.46
1:B:172:GLU:HG3	1:B:178:LYS:CG	2.43	0.46
1:C:216:PRO:HA	1:C:233:LYS:HE2	1.96	0.46
1:A:54:GLN:HG2	1:A:59:ASN:OD1	2.16	0.45
1:B:23:MSE:CE	1:B:46:LEU:HD11	2.46	0.45
1:C:103:TYR:CE1	1:C:130:ILE:CD1	2.98	0.45
1:C:204:MSE:HB3	1:C:207:ILE:HG22	1.98	0.45
1:A:1:MSE:HE1	1:A:248:LEU:HD21	1.97	0.45
1:A:90:ASP:HA	1:A:93:ASN:ND2	2.32	0.45
1:A:124:ALA:CB	1:A:159:ASP:HB2	2.46	0.45
1:C:49:HIS:CB	1:C:273:ARG:NH2	2.79	0.45
1:A:54:GLN:NE2	1:A:96:LEU:HD11	2.32	0.45
1:A:191:ILE:HD13	1:A:236:MSE:SE	2.66	0.45
1:C:177:ILE:HD12	1:C:194:ILE:HG12	1.99	0.45
1:A:134:GLN:HE21	1:A:136:ASP:H	1.60	0.45
1:A:134:GLN:HE21	1:A:136:ASP:N	2.15	0.45
1:A:260:TYR:CA	1:A:267:ILE:HD12	2.46	0.45
1:B:2:LYS:NZ	1:B:2:LYS:HB2	2.31	0.45
1:C:101:ILE:HG12	1:C:126:PHE:CD2	2.51	0.45
1:A:72:SER:O	1:A:76:LEU:HG	2.18	0.44
1:B:29:LYS:HG3	1:B:32:TRP:CZ2	2.52	0.44
1:B:148:SER:HB3	1:B:152:TYR:HD2	1.80	0.44
1:C:207:ILE:HD12	1:C:207:ILE:HA	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ASN:OD1	1:B:37:ASN:N	2.51	0.44
1:A:260:TYR:HA	1:A:267:ILE:HA	1.99	0.44
1:C:60:LEU:HD12	1:C:105:LEU:HD11	1.98	0.44
1:C:119:ASP:C	1:C:119:ASP:OD2	2.60	0.44
1:B:52:LEU:HD13	1:B:61:ILE:HD11	2.00	0.44
1:B:29:LYS:C	1:B:31:LEU:N	2.74	0.44
1:B:274:GLU:O	1:B:275:THR:C	2.61	0.44
1:A:17:LYS:HB2	1:A:80:GLY:HA3	2.00	0.43
1:A:267:ILE:O	1:A:267:ILE:HG13	2.18	0.43
1:B:39:ASN:OD1	1:B:39:ASN:C	2.61	0.43
1:B:134:GLN:HE22	1:B:136:ASP:CB	2.31	0.43
1:C:170:HIS:HA	1:C:179:MSE:O	2.18	0.43
1:A:216:PRO:HA	1:A:233:LYS:HE2	1.98	0.43
1:C:17:LYS:O	1:C:18:MSE:HE2	2.17	0.43
1:A:1:MSE:CE	1:A:3:ILE:HD11	2.49	0.43
1:A:87:ILE:HG23	1:A:88:ILE:HD12	2.00	0.43
1:B:134:GLN:HG2	1:B:188:GLY:O	2.18	0.43
1:C:16:THR:HB	1:C:18:MSE:CE	2.48	0.43
1:C:62:ILE:O	1:C:63:ASP:HB2	2.18	0.43
1:C:123:HIS:CD2	1:C:159:ASP:OD2	2.72	0.43
1:A:131:HIS:CE1	1:A:160:TYR:HB2	2.53	0.43
1:B:70:LYS:HZ3	1:B:156:ASN:HD21	1.65	0.43
1:B:87:ILE:HG23	1:B:88:ILE:N	2.33	0.43
1:C:177:ILE:CD1	1:C:194:ILE:HG23	2.48	0.43
1:C:53:ILE:CD1	1:C:194:ILE:HD11	2.48	0.43
1:C:210:THR:HB	1:C:252:ASP:OD1	2.17	0.43
1:A:3:ILE:O	1:A:196:SER:HB3	2.18	0.43
1:B:29:LYS:N	1:B:30:PRO:CD	2.82	0.43
1:A:259:LYS:N	1:A:267:ILE:HD11	2.33	0.43
1:B:66:ILE:HA	1:B:115:ALA:HB3	1.99	0.43
1:C:127:GLU:O	1:C:128:ASN:CB	2.66	0.43
1:C:27:VAL:HG21	1:C:218:TRP:CD2	2.54	0.43
1:A:88:ILE:HG13	1:A:98:PRO:HG2	2.00	0.43
1:A:191:ILE:HD13	1:A:236:MSE:HE1	2.00	0.43
1:A:195:GLU:HG2	1:A:200:LYS:HG2	2.01	0.43
1:A:267:ILE:CG2	1:A:270:TYR:CE2	3.02	0.43
1:B:137:GLU:HB2	1:B:188:GLY:HA3	2.00	0.43
1:B:106:MSE:HE1	1:B:113:HIS:O	2.19	0.42
1:B:204:MSE:HB3	1:B:207:ILE:HG22	2.00	0.42
1:B:252:ASP:O	1:B:273:ARG:NH1	2.52	0.42
1:A:97:THR:C	1:A:99:LYS:N	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:134:GLN:NE2	1:C:136:ASP:N	2.65	0.42
1:B:112:ASP:OD2	1:B:113:HIS:CE1	2.73	0.42
1:C:101:ILE:HG13	1:C:126:PHE:CD2	2.54	0.42
1:C:252:ASP:O	1:C:273:ARG:HD2	2.19	0.42
1:A:87:ILE:CG2	1:A:88:ILE:HD12	2.49	0.42
1:B:105:LEU:HD23	1:B:132:VAL:HB	2.01	0.42
1:A:259:LYS:HB3	1:A:269:ALA:HB3	2.02	0.42
1:A:267:ILE:CG2	1:A:270:TYR:CZ	3.02	0.42
1:B:103:TYR:CD1	1:B:130:ILE:HB	2.54	0.42
1:A:197:GLN:HB3	1:A:198:GLY:H	1.55	0.41
1:B:170:HIS:HA	1:B:179:MSE:O	2.20	0.41
1:C:72:SER:OG	1:C:75:GLN:HG3	2.20	0.41
1:C:97:THR:CB	1:C:98:PRO:HD2	2.49	0.41
1:A:236:MSE:HB3	1:A:240:PHE:CE2	2.55	0.41
1:C:118:THR:HG21	1:C:158:GLY:HA3	2.02	0.41
1:C:261:SER:C	1:C:263:ASP:H	2.28	0.41
1:A:1:MSE:HB2	1:A:8:ILE:O	2.21	0.41
1:A:62:ILE:O	1:A:62:ILE:HG22	2.19	0.41
1:B:197:GLN:HB3	1:B:198:GLY:H	1.68	0.41
1:C:100:ASP:O	1:C:101:ILE:O	2.39	0.41
1:C:41:ARG:HG3	1:C:41:ARG:HH11	1.85	0.41
1:C:191:ILE:HG22	1:C:204:MSE:HA	2.03	0.41
1:C:2:LYS:O	1:C:2:LYS:HG2	2.20	0.41
1:B:270:TYR:CD1	1:B:270:TYR:C	2.99	0.41
1:C:193:THR:HG22	1:C:202:VAL:CG2	2.50	0.41
1:C:273:ARG:C	1:C:274:GLU:HG3	2.45	0.41
1:A:246:TRP:CE2	1:A:259:LYS:HE2	2.56	0.41
1:A:20:GLY:HA3	1:A:32:TRP:CH2	2.56	0.41
1:A:263:ASP:HB3	1:A:265:GLU:OE2	2.21	0.41
1:B:67:GLY:HA2	2:B:309:HOH:O	2.19	0.41
1:B:128:ASN:H	1:B:163:LYS:NZ	2.16	0.41
1:C:27:VAL:HG21	1:C:218:TRP:CG	2.56	0.41
1:C:123:HIS:CE1	1:C:127:GLU:OE1	2.61	0.41
1:C:148:SER:HB3	1:C:152:TYR:CD2	2.52	0.41
1:A:267:ILE:O	1:A:270:TYR:CE2	2.74	0.41
1:C:50:PRO:CG	1:C:65:GLY:H	2.29	0.41
1:A:51:ILE:HB	1:A:62:ILE:CG1	2.50	0.40
1:B:24:PHE:HB3	1:B:27:VAL:HB	2.02	0.40
1:B:127:GLU:H	1:B:127:GLU:HG2	1.73	0.40
1:C:98:PRO:O	1:C:99:LYS:HB2	2.20	0.40
1:A:92:ALA:C	1:A:94:TYR:N	2.80	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:THR:HG22	1:A:202:VAL:HG23	2.03	0.40
1:C:97:THR:CG2	1:C:98:PRO:N	2.82	0.40
1:C:200:LYS:HD3	1:C:245:TYR:CE1	2.57	0.40
1:A:63:ASP:HB2	1:A:250:TYR:CE1	2.56	0.40
1:C:87:ILE:HG23	1:C:88:ILE:H	1.85	0.40
1:C:101:ILE:HG13	1:C:126:PHE:CE2	2.57	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	275/280 (98%)	250 (91%)	18 (6%)	7 (2%)	4	7
1	B	273/280 (98%)	252 (92%)	18 (7%)	3 (1%)	11	22
1	C	273/280 (98%)	243 (89%)	23 (8%)	7 (3%)	4	7
All	All	821/840 (98%)	745 (91%)	59 (7%)	17 (2%)	5	9

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	HIS
1	A	124	ALA
1	A	262	ASP
1	B	39	ASN
1	B	123	HIS
1	B	124	ALA
1	C	99	LYS
1	C	60	LEU
1	C	101	ILE
1	C	274	GLU
1	C	262	ASP

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Mol	Chain	Res	Type
1	A	99	LYS
1	C	98	PRO
1	A	159	ASP
1	A	263	ASP
1	A	173	PRO
1	C	65	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	240/234 (103%)	229 (95%)	11 (5%)	24	48
1	B	238/234 (102%)	226 (95%)	12 (5%)	22	44
1	C	238/234 (102%)	223 (94%)	15 (6%)	16	34
All	All	716/702 (102%)	678 (95%)	38 (5%)	20	42

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	42	ASN
1	A	60	LEU
1	A	73	GLU
1	A	90	ASP
1	A	134	GLN
1	A	145	ASN
1	A	197	GLN
1	A	207	ILE
1	A	235	ARG
1	A	274	GLU
1	B	1	MSE
1	B	2	LYS
1	B	40	GLU
1	B	41	ARG
1	B	60	LEU

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Mol	Chain	Res	Type
1	B	134	GLN
1	B	145	ASN
1	B	149	LYS
1	B	159	ASP
1	B	197	GLN
1	B	207	ILE
1	B	235	ARG
1	C	2	LYS
1	C	5	ASP
1	C	59	ASN
1	C	60	LEU
1	C	97	THR
1	C	134	GLN
1	C	145	ASN
1	C	148	SER
1	C	149	LYS
1	C	151	THR
1	C	169	LYS
1	C	197	GLN
1	C	207	ILE
1	C	235	ARG
1	C	263	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	54	GLN
1	A	57	GLN
1	A	93	ASN
1	A	113	HIS
1	A	128	ASN
1	A	131	HIS
1	A	134	GLN
1	A	135	GLN
1	A	145	ASN
1	A	156	ASN
1	A	180	GLN
1	A	181	HIS
1	A	197	GLN
1	A	244	GLN
1	A	254	ASN

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Mol	Chain	Res	Type
1	B	42	ASN
1	B	49	HIS
1	B	54	GLN
1	B	113	HIS
1	B	131	HIS
1	B	134	GLN
1	B	145	ASN
1	B	156	ASN
1	B	197	GLN
1	B	244	GLN
1	C	9	HIS
1	C	42	ASN
1	C	68	ASN
1	C	110	HIS
1	C	113	HIS
1	C	123	HIS
1	C	128	ASN
1	C	131	HIS
1	C	134	GLN
1	C	135	GLN
1	C	145	ASN
1	C	156	ASN
1	C	180	GLN
1	C	181	HIS
1	C	197	GLN
1	C	244	GLN
1	C	254	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	268/280 (95%)	0.75	25 (9%) 14 12	25, 42, 63, 86	1 (0%)
1	B	266/280 (95%)	0.82	26 (9%) 13 11	25, 49, 67, 77	1 (0%)
1	C	266/280 (95%)	0.88	31 (11%) 9 7	23, 47, 66, 79	0
All	All	800/840 (95%)	0.82	82 (10%) 12 10	23, 45, 66, 86	2 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	267	ILE	10.5
1	C	98	PRO	6.5
1	A	262	ASP	6.2
1	A	265	GLU	5.9
1	A	263	ASP	5.7
1	A	124	ALA	5.2
1	C	124	ALA	5.0
1	C	59	ASN	5.0
1	C	103	TYR	4.8
1	C	275	THR	4.8
1	C	274	GLU	4.5
1	A	264	GLY	4.5
1	C	97	THR	4.3
1	A	98	PRO	4.2
1	C	64	ALA	4.2
1	C	95	ASN	4.0
1	A	125	ILE	4.0
1	A	155	LYS	3.9
1	A	123	HIS	3.8
1	B	262	ASP	3.8
1	C	63	ASP	3.8
1	C	102	ASP	3.8
1	B	267	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	93	ASN	3.7
1	A	97	THR	3.6
1	C	2	LYS	3.6
1	A	121	ALA	3.5
1	C	262	ASP	3.5
1	B	37	ASN	3.5
1	B	38	ALA	3.4
1	B	155	LYS	3.3
1	A	126	PHE	3.3
1	A	266	ASN	3.3
1	B	36	TYR	3.3
1	B	270	TYR	3.2
1	A	277	VAL	3.2
1	C	100	ASP	3.2
1	C	62	ILE	3.1
1	B	275	THR	3.1
1	B	95	ASN	2.9
1	C	101	ILE	2.9
1	C	65	GLY	2.9
1	B	269	ALA	2.9
1	B	124	ALA	2.8
1	B	125	ILE	2.7
1	A	83	GLU	2.7
1	C	272	LEU	2.6
1	C	80	GLY	2.6
1	B	266	ASN	2.6
1	B	126	PHE	2.6
1	A	96	LEU	2.5
1	C	195	GLU	2.4
1	A	92	ALA	2.4
1	A	122	GLY	2.4
1	C	66	ILE	2.3
1	A	268	ASP	2.3
1	B	73	GLU	2.3
1	B	254	ASN	2.3
1	C	207	ILE	2.3
1	C	265	GLU	2.3
1	B	91	LEU	2.3
1	B	268	ASP	2.3
1	C	263	ASP	2.3
1	B	10	TYR	2.3
1	A	197	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	272	LEU	2.2
1	C	60	LEU	2.2
1	C	61	ILE	2.2
1	A	40	GLU	2.2
1	B	40	GLU	2.2
1	B	30	PRO	2.2
1	C	125	ILE	2.1
1	B	127	GLU	2.1
1	C	94	TYR	2.1
1	B	263	ASP	2.1
1	A	276	LEU	2.1
1	B	39	ASN	2.1
1	C	116	GLY	2.1
1	A	274	GLU	2.1
1	B	33	SER	2.0
1	C	96	LEU	2.0
1	C	5	ASP	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.