



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:51 PM UTC

PDB ID : 3GUY / pdb_00003guy
Title : Crystal structure of a short-chain dehydrogenase/reductase from *Vibrio parahaemolyticus*
Authors : Patskovsky, Y.; Bonanno, J.B.; Freeman, J.; Bain, K.T.; Miller, S.; Sampathkumar, P.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-03-30
Resolution : 1.90 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
Xtrriage (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

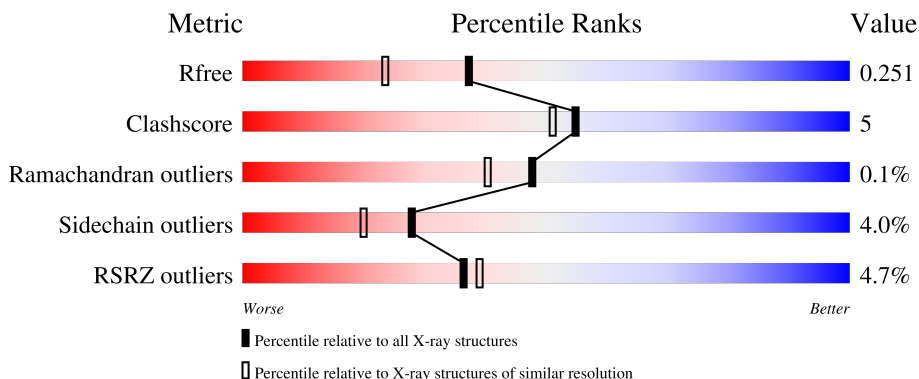
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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.



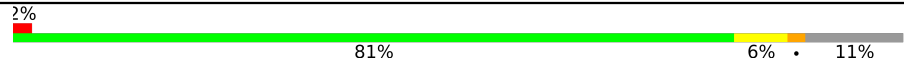

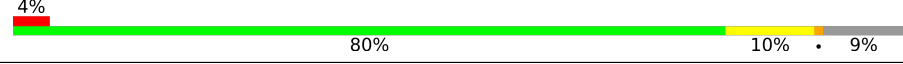
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	230	2% 82% 7% 11%
1	B	230	3% 81% 7% 11%
1	C	230	4% 84% 7% 8%
1	D	230	5% 86% 9% 8%
1	E	230	2% 79% 10% 11%

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Mol	Chain	Length	Quality of chain
1	F	230	 <p>2% 81% 6% • 11%</p>
1	G	230	 <p>12% 81% 15% • •</p>
1	H	230	 <p>4% 80% 10% • 9%</p>

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 13319 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 Short-chain dehydrogenase/reductase SDR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	204	Total 1539	C 964	N 264	O 303	S 8	0	0	0
1	B	204	Total 1540	C 965	N 264	O 303	S 8	0	0	0
1	C	212	Total 1598	C 1003	N 273	O 314	S 8	0	0	0
1	D	220	Total 1650	C 1034	N 281	O 327	S 8	0	0	0
1	E	204	Total 1545	C 970	N 264	O 303	S 8	0	0	0
1	F	204	Total 1541	C 967	N 263	O 303	S 8	0	0	0
1	G	228	Total 1735	C 1082	N 302	O 343	S 8	0	0	0
1	H	209	Total 1585	C 999	N 267	O 311	S 8	0	0	0

There are 104 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP A6B7Q2
A	0	SER	-	expression tag	UNP A6B7Q2
A	1	LEU	-	expression tag	UNP A6B7Q2
A	52	ARG	CYS	engineered mutation	UNP A6B7Q2
A	204	ALA	SER	engineered mutation	UNP A6B7Q2
A	221	GLU	-	expression tag	UNP A6B7Q2
A	222	GLY	-	expression tag	UNP A6B7Q2
A	223	HIS	-	expression tag	UNP A6B7Q2
A	224	HIS	-	expression tag	UNP A6B7Q2
A	225	HIS	-	expression tag	UNP A6B7Q2
A	226	HIS	-	expression tag	UNP A6B7Q2
A	227	HIS	-	expression tag	UNP A6B7Q2
A	228	HIS	-	expression tag	UNP A6B7Q2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MET	-	expression tag	UNP A6B7Q2
B	0	SER	-	expression tag	UNP A6B7Q2
B	1	LEU	-	expression tag	UNP A6B7Q2
B	52	ARG	CYS	engineered mutation	UNP A6B7Q2
B	204	ALA	SER	engineered mutation	UNP A6B7Q2
B	221	GLU	-	expression tag	UNP A6B7Q2
B	222	GLY	-	expression tag	UNP A6B7Q2
B	223	HIS	-	expression tag	UNP A6B7Q2
B	224	HIS	-	expression tag	UNP A6B7Q2
B	225	HIS	-	expression tag	UNP A6B7Q2
B	226	HIS	-	expression tag	UNP A6B7Q2
B	227	HIS	-	expression tag	UNP A6B7Q2
B	228	HIS	-	expression tag	UNP A6B7Q2
C	-1	MET	-	expression tag	UNP A6B7Q2
C	0	SER	-	expression tag	UNP A6B7Q2
C	1	LEU	-	expression tag	UNP A6B7Q2
C	52	ARG	CYS	engineered mutation	UNP A6B7Q2
C	204	ALA	SER	engineered mutation	UNP A6B7Q2
C	221	GLU	-	expression tag	UNP A6B7Q2
C	222	GLY	-	expression tag	UNP A6B7Q2
C	223	HIS	-	expression tag	UNP A6B7Q2
C	224	HIS	-	expression tag	UNP A6B7Q2
C	225	HIS	-	expression tag	UNP A6B7Q2
C	226	HIS	-	expression tag	UNP A6B7Q2
C	227	HIS	-	expression tag	UNP A6B7Q2
C	228	HIS	-	expression tag	UNP A6B7Q2
D	-1	MET	-	expression tag	UNP A6B7Q2
D	0	SER	-	expression tag	UNP A6B7Q2
D	1	LEU	-	expression tag	UNP A6B7Q2
D	52	ARG	CYS	engineered mutation	UNP A6B7Q2
D	204	ALA	SER	engineered mutation	UNP A6B7Q2
D	221	GLU	-	expression tag	UNP A6B7Q2
D	222	GLY	-	expression tag	UNP A6B7Q2
D	223	HIS	-	expression tag	UNP A6B7Q2
D	224	HIS	-	expression tag	UNP A6B7Q2
D	225	HIS	-	expression tag	UNP A6B7Q2
D	226	HIS	-	expression tag	UNP A6B7Q2
D	227	HIS	-	expression tag	UNP A6B7Q2
D	228	HIS	-	expression tag	UNP A6B7Q2
E	-1	MET	-	expression tag	UNP A6B7Q2
E	0	SER	-	expression tag	UNP A6B7Q2
E	1	LEU	-	expression tag	UNP A6B7Q2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	52	ARG	CYS	engineered mutation	UNP A6B7Q2
E	204	ALA	SER	engineered mutation	UNP A6B7Q2
E	221	GLU	-	expression tag	UNP A6B7Q2
E	222	GLY	-	expression tag	UNP A6B7Q2
E	223	HIS	-	expression tag	UNP A6B7Q2
E	224	HIS	-	expression tag	UNP A6B7Q2
E	225	HIS	-	expression tag	UNP A6B7Q2
E	226	HIS	-	expression tag	UNP A6B7Q2
E	227	HIS	-	expression tag	UNP A6B7Q2
E	228	HIS	-	expression tag	UNP A6B7Q2
F	-1	MET	-	expression tag	UNP A6B7Q2
F	0	SER	-	expression tag	UNP A6B7Q2
F	1	LEU	-	expression tag	UNP A6B7Q2
F	52	ARG	CYS	engineered mutation	UNP A6B7Q2
F	204	ALA	SER	engineered mutation	UNP A6B7Q2
F	221	GLU	-	expression tag	UNP A6B7Q2
F	222	GLY	-	expression tag	UNP A6B7Q2
F	223	HIS	-	expression tag	UNP A6B7Q2
F	224	HIS	-	expression tag	UNP A6B7Q2
F	225	HIS	-	expression tag	UNP A6B7Q2
F	226	HIS	-	expression tag	UNP A6B7Q2
F	227	HIS	-	expression tag	UNP A6B7Q2
F	228	HIS	-	expression tag	UNP A6B7Q2
G	-1	MET	-	expression tag	UNP A6B7Q2
G	0	SER	-	expression tag	UNP A6B7Q2
G	1	LEU	-	expression tag	UNP A6B7Q2
G	52	ARG	CYS	engineered mutation	UNP A6B7Q2
G	204	ALA	SER	engineered mutation	UNP A6B7Q2
G	221	GLU	-	expression tag	UNP A6B7Q2
G	222	GLY	-	expression tag	UNP A6B7Q2
G	223	HIS	-	expression tag	UNP A6B7Q2
G	224	HIS	-	expression tag	UNP A6B7Q2
G	225	HIS	-	expression tag	UNP A6B7Q2
G	226	HIS	-	expression tag	UNP A6B7Q2
G	227	HIS	-	expression tag	UNP A6B7Q2
G	228	HIS	-	expression tag	UNP A6B7Q2
H	-1	MET	-	expression tag	UNP A6B7Q2
H	0	SER	-	expression tag	UNP A6B7Q2
H	1	LEU	-	expression tag	UNP A6B7Q2
H	52	ARG	CYS	engineered mutation	UNP A6B7Q2
H	204	ALA	SER	engineered mutation	UNP A6B7Q2
H	221	GLU	-	expression tag	UNP A6B7Q2

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Chain	Residue	Modelled	Actual	Comment	Reference
H	222	GLY	-	expression tag	UNP A6B7Q2
H	223	HIS	-	expression tag	UNP A6B7Q2
H	224	HIS	-	expression tag	UNP A6B7Q2
H	225	HIS	-	expression tag	UNP A6B7Q2
H	226	HIS	-	expression tag	UNP A6B7Q2
H	227	HIS	-	expression tag	UNP A6B7Q2
H	228	HIS	-	expression tag	UNP A6B7Q2

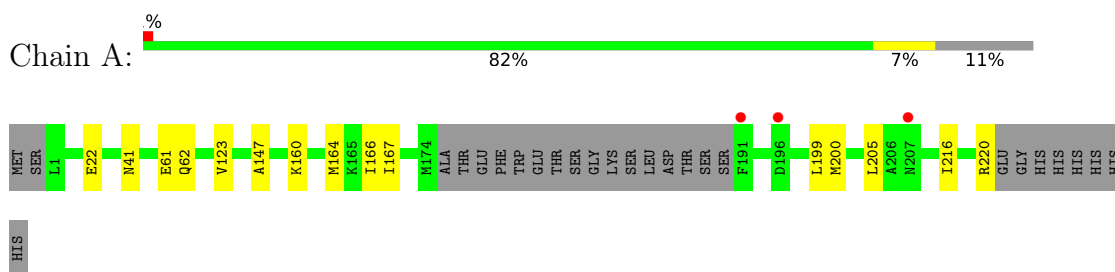
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	86	Total O 86 86	0	0
2	B	92	Total O 92 92	0	0
2	C	91	Total O 91 91	0	0
2	D	74	Total O 74 74	0	0
2	E	63	Total O 63 63	0	0
2	F	66	Total O 66 66	0	0
2	G	61	Total O 61 61	0	0
2	H	53	Total O 53 53	0	0

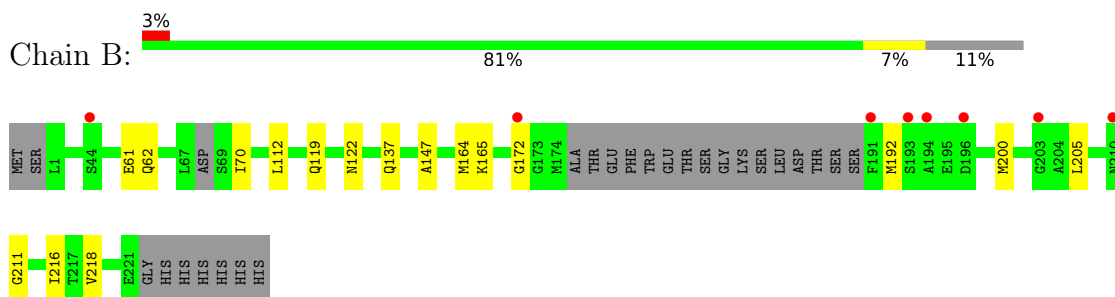
3 Residue-property plots [i](#)

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.

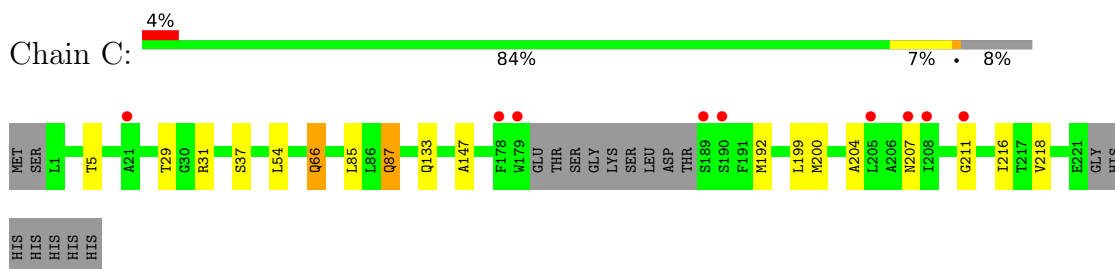
- Molecule 1: Short-chain dehydrogenase/reductase SDR



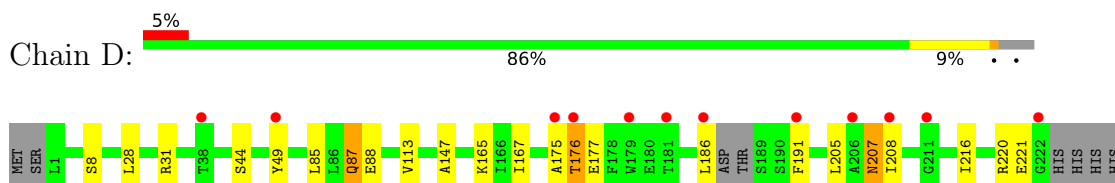
- Molecule 1: Short-chain dehydrogenase/reductase SDR



- Molecule 1: Short-chain dehydrogenase/reductase SDR



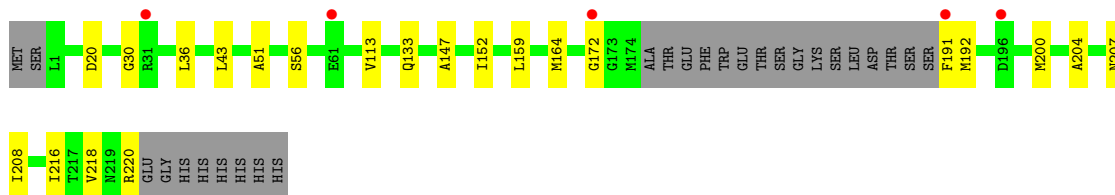
- Molecule 1: Short-chain dehydrogenase/reductase SDR



HIS
HIS

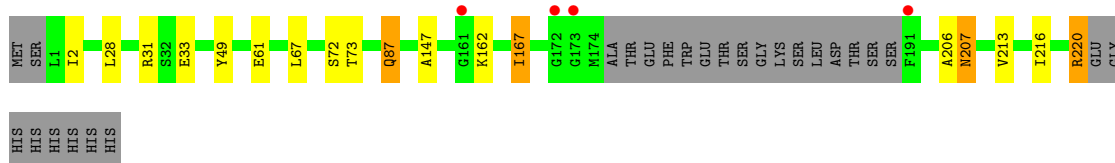
• Molecule 1: Short-chain dehydrogenase/reductase SDR

Chain E: 2% 79% 10% 11%



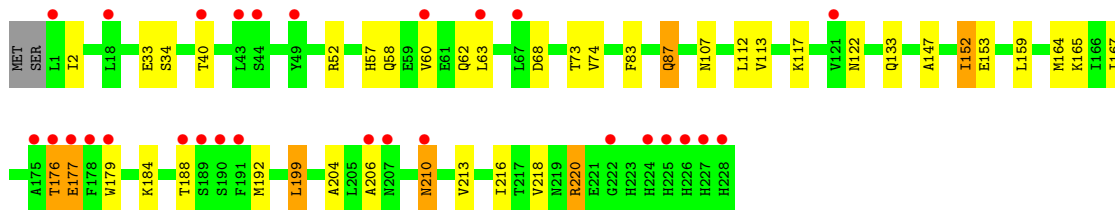
• Molecule 1: Short-chain dehydrogenase/reductase SDR

Chain F: 2% 81% 6% 11%



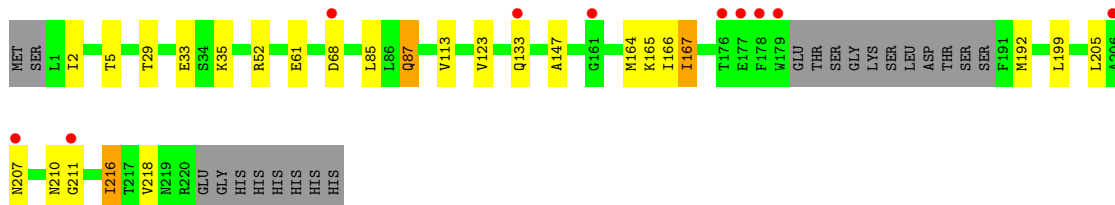
• Molecule 1: Short-chain dehydrogenase/reductase SDR

Chain G: 12% 81% 15% 2%



• Molecule 1: Short-chain dehydrogenase/reductase SDR

Chain H: 4% 80% 10% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.81Å 71.31Å 93.08Å 72.38° 89.39° 74.30°	Depositor
Resolution (Å)	20.00 – 1.90 20.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.9 (20.00-1.90) 96.2 (20.00-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 1.89Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.200 , 0.249 0.205 , 0.251	Depositor DCC
R_{free} test set	6220 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtrriage
Anisotropy	0.121	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13319	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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	1.14	1/1558 (0.1%)	0.99	0/2107
1	B	1.06	0/1558	0.97	0/2105
1	C	1.05	0/1620	0.99	0/2193
1	D	1.05	0/1672	0.98	0/2261
1	E	0.94	0/1565	0.91	0/2116
1	F	1.00	1/1561 (0.1%)	0.95	0/2112
1	G	0.91	0/1764	0.99	1/2387 (0.0%)
1	H	0.92	0/1608	0.97	1/2177 (0.0%)
All	All	1.01	2/12906 (0.0%)	0.97	2/17458 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	213	VAL	CA-CB	6.26	1.60	1.53
1	A	62	GLN	N-CA	5.01	1.52	1.46

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	211	GLY	N-CA-C	5.78	119.58	110.96
1	G	40	THR	N-CA-C	5.06	117.53	111.71

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1539	0	1559	8	0
1	B	1540	0	1560	17	0
1	C	1598	0	1604	19	0
1	D	1650	0	1652	18	0
1	E	1545	0	1566	20	0
1	F	1541	0	1555	10	0
1	G	1735	0	1719	30	0
1	H	1585	0	1592	19	0
2	A	86	0	0	1	0
2	B	92	0	0	0	0
2	C	91	0	0	3	0
2	D	74	0	0	1	0
2	E	63	0	0	1	0
2	F	66	0	0	1	0
2	G	61	0	0	1	0
2	H	53	0	0	0	0
All	All	13319	0	12807	123	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 5.

All (123) 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:205:LEU:HD21	1:B:216:ILE:HD12	1.39	1.02
1:B:205:LEU:CD2	1:B:216:ILE:CD1	2.36	1.02
1:B:205:LEU:HD23	1:B:216:ILE:HD11	1.44	0.99
1:B:205:LEU:CD2	1:B:216:ILE:HD11	1.93	0.96
1:E:159:LEU:HD11	1:E:164:MET:HE2	1.47	0.95
1:D:87:GLN:HE21	1:D:87:GLN:H	1.16	0.90
1:B:205:LEU:HD23	1:B:216:ILE:CD1	2.00	0.87
1:F:87:GLN:HE21	1:F:87:GLN:H	1.23	0.86
1:G:147:ALA:HB2	1:H:147:ALA:HB2	1.59	0.84
1:B:205:LEU:HD21	1:B:216:ILE:CD1	2.02	0.84
1:D:167:ILE:HG21	1:D:205:LEU:HD21	1.59	0.83
1:E:204:ALA:HB1	1:E:216:ILE:CD1	2.09	0.82
1:G:113:VAL:HA	1:G:164:MET:HE1	1.68	0.76
1:E:204:ALA:CB	1:E:216:ILE:CD1	2.64	0.75
1:H:192:MET:SD	1:H:218:VAL:HG12	2.29	0.73
1:H:167:ILE:HG21	1:H:205:LEU:HD21	1.74	0.70
1:H:167:ILE:HG21	1:H:205:LEU:CD2	2.22	0.70
1:C:147:ALA:HB2	1:D:147:ALA:HB2	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ASN:HB3	2:A:542:HOH:O	1.96	0.66
1:A:123:VAL:HG21	1:A:164:MET:HE3	1.77	0.66
1:G:117:LYS:HA	1:G:164:MET:HE2	1.77	0.65
1:E:204:ALA:CB	1:E:216:ILE:HD13	2.26	0.65
1:E:147:ALA:HB2	1:F:147:ALA:HB2	1.78	0.64
1:D:167:ILE:HG21	1:D:205:LEU:CD2	2.28	0.64
1:H:85:LEU:HD13	1:H:87:GLN:NE2	2.13	0.63
1:E:207:ASN:O	1:E:208:ILE:HD13	1.99	0.63
1:D:31:ARG:NH1	2:D:721:HOH:O	2.31	0.63
1:A:147:ALA:HB2	1:B:147:ALA:HB2	1.81	0.63
1:B:211:GLY:C	1:C:200:MET:HE1	2.23	0.63
1:H:123:VAL:HG21	1:H:164:MET:HE3	1.80	0.62
1:G:192:MET:HB2	1:G:220:ARG:HB2	1.84	0.60
1:E:204:ALA:HB3	1:E:216:ILE:HD13	1.84	0.59
1:G:87:GLN:HG2	1:H:113:VAL:CG1	2.32	0.59
1:E:159:LEU:CD1	1:E:164:MET:HE2	2.28	0.59
1:F:206:ALA:C	1:F:207:ASN:HD22	2.10	0.59
1:B:70:ILE:HD13	1:B:119:GLN:HG3	1.85	0.58
1:D:87:GLN:H	1:D:87:GLN:NE2	1.95	0.58
1:C:31:ARG:NH2	2:C:694:HOH:O	2.35	0.58
1:H:166:ILE:C	1:H:167:ILE:HD13	2.29	0.58
1:G:60:VAL:HG21	1:G:107:ASN:HB3	1.85	0.58
1:E:113:VAL:CG2	1:E:164:MET:HE1	2.34	0.57
1:B:172:GLY:O	1:B:192:MET:HE2	2.04	0.57
1:E:113:VAL:HG22	1:E:164:MET:HE1	1.85	0.56
1:G:113:VAL:HG22	1:G:164:MET:HE1	1.88	0.55
1:E:172:GLY:O	1:E:192:MET:HE2	2.07	0.55
1:C:87:GLN:HG3	1:D:113:VAL:HG12	1.89	0.55
1:B:122:ASN:OD1	1:B:165:LYS:NZ	2.40	0.54
1:C:192:MET:SD	1:C:218:VAL:HG12	2.47	0.54
1:D:175:ALA:O	1:D:176:THR:C	2.51	0.53
1:B:200:MET:HE1	1:C:211:GLY:C	2.33	0.53
1:C:87:GLN:NE2	2:C:539:HOH:O	2.41	0.53
1:B:192:MET:SD	1:B:218:VAL:HG12	2.48	0.53
1:F:207:ASN:HD22	1:F:207:ASN:N	2.07	0.53
1:F:87:GLN:H	1:F:87:GLN:NE2	2.02	0.52
1:H:192:MET:SD	1:H:218:VAL:CG1	2.98	0.52
1:H:85:LEU:HD13	1:H:87:GLN:HE21	1.74	0.51
1:E:192:MET:SD	1:E:218:VAL:HG12	2.50	0.51
1:C:85:LEU:HB3	1:C:87:GLN:HE22	1.75	0.51
1:G:152:ILE:HD12	1:G:152:ILE:C	2.34	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:191:PHE:HA	1:E:220:ARG:HD3	1.92	0.51
1:H:167:ILE:HD13	1:H:167:ILE:N	2.25	0.51
1:E:208:ILE:HG21	1:G:199:LEU:HD12	1.92	0.51
1:G:57:HIS:CE1	1:G:58:GLN:OE1	2.64	0.51
1:B:200:MET:HE1	1:C:211:GLY:HA3	1.94	0.50
1:G:159:LEU:HD11	1:G:164:MET:HE3	1.93	0.50
1:G:204:ALA:CB	1:G:216:ILE:CD1	2.89	0.50
1:G:87:GLN:HG2	1:H:113:VAL:HG11	1.94	0.49
1:G:87:GLN:HG2	1:H:113:VAL:HG12	1.93	0.49
1:C:204:ALA:CB	1:C:216:ILE:CD1	2.89	0.49
1:C:87:GLN:CG	1:D:113:VAL:HG12	2.42	0.49
1:F:220:ARG:C	1:H:210:ASN:O	2.56	0.48
1:E:200:MET:HG2	1:G:206:ALA:HB3	1.95	0.48
1:G:152:ILE:HD12	1:G:153:GLU:N	2.28	0.48
1:F:31:ARG:NH1	2:F:741:HOH:O	2.46	0.48
1:A:167:ILE:HG21	1:A:205:LEU:HD22	1.96	0.47
1:F:167:ILE:HD13	1:F:167:ILE:N	2.30	0.47
1:G:2:ILE:HD12	1:G:73:THR:HB	1.97	0.47
1:D:85:LEU:HB2	1:D:88:GLU:HG2	1.97	0.47
1:A:166:ILE:C	1:A:167:ILE:HD13	2.39	0.46
2:E:329:HOH:O	1:G:210:ASN:HB2	2.14	0.46
1:G:52:ARG:HG3	1:G:63:LEU:HD13	1.97	0.46
1:B:112:LEU:O	1:B:164:MET:HE1	2.15	0.46
1:G:113:VAL:HG22	1:G:164:MET:CE	2.45	0.46
1:G:176:THR:HG22	1:G:177:GLU:N	2.31	0.45
1:C:204:ALA:CB	1:C:216:ILE:HD13	2.46	0.45
1:D:28:LEU:O	1:D:49:TYR:HA	2.17	0.45
1:E:204:ALA:HB1	1:E:216:ILE:HD11	1.93	0.45
1:H:5:THR:HA	1:H:29:THR:HG22	1.97	0.45
1:H:52:ARG:HA	1:H:52:ARG:HD3	1.86	0.44
1:A:216:ILE:HG22	1:D:216:ILE:HG13	1.99	0.44
1:E:30:GLY:O	1:E:51:ALA:HA	2.17	0.44
1:H:85:LEU:HB3	1:H:87:GLN:HE21	1.82	0.44
1:G:192:MET:SD	1:G:218:VAL:HG12	2.58	0.44
1:E:30:GLY:N	1:E:36:LEU:HD11	2.32	0.44
1:C:31:ARG:NH1	2:C:450:HOH:O	2.40	0.44
1:G:176:THR:CG2	1:G:177:GLU:N	2.80	0.44
1:H:165:LYS:NZ	1:H:207:ASN:OD1	2.36	0.43
1:F:2:ILE:HD12	1:F:73:THR:HB	1.99	0.43
1:F:28:LEU:O	1:F:49:TYR:HA	2.18	0.43
1:G:133:GLN:HB3	2:G:239:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:74:VAL:HG11	1:G:112:LEU:HD11	2.01	0.43
1:D:175:ALA:O	1:D:177:GLU:N	2.52	0.43
1:D:186:LEU:HD22	1:D:191:PHE:CZ	2.54	0.43
1:A:167:ILE:HD13	1:A:167:ILE:N	2.33	0.43
1:E:20:ASP:CB	1:E:43:LEU:HD22	2.49	0.43
1:D:208:ILE:HG22	1:D:208:ILE:O	2.18	0.42
1:G:204:ALA:HB1	1:G:216:ILE:CD1	2.49	0.42
1:G:122:ASN:OD1	1:G:165:LYS:CE	2.67	0.42
1:C:5:THR:HB	1:C:54:LEU:HD11	2.01	0.42
1:G:204:ALA:O	1:G:213:VAL:HG21	2.20	0.42
1:H:216:ILE:HD13	1:H:216:ILE:HG21	1.67	0.42
1:G:83:PHE:HB2	1:G:179:TRP:CE3	2.54	0.42
1:A:200:MET:SD	1:D:207:ASN:HB2	2.60	0.41
1:C:5:THR:HA	1:C:29:THR:HG22	2.02	0.41
1:D:85:LEU:HB2	1:D:88:GLU:CG	2.50	0.41
1:B:61:GLU:OE1	1:B:62:GLN:NE2	2.53	0.41
1:E:172:GLY:O	1:E:192:MET:CE	2.68	0.41
1:C:204:ALA:HB1	1:C:216:ILE:CD1	2.50	0.41
1:C:66:GLN:CA	1:C:66:GLN:OE1	2.69	0.41
1:B:200:MET:HE1	1:C:211:GLY:CA	2.52	0.40
1:D:87:GLN:HE21	1:D:87:GLN:N	1.98	0.40
1:G:87:GLN:H	1:G:87:GLN:HG3	1.43	0.40
1:C:87:GLN:NE2	1:C:87:GLN:H	2.19	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	200/230 (87%)	197 (98%)	3 (2%)	0	100 100
1	B	198/230 (86%)	195 (98%)	3 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	208/230 (90%)	204 (98%)	4 (2%)	0	100	100
1	D	216/230 (94%)	212 (98%)	3 (1%)	1 (0%)	24	16
1	E	200/230 (87%)	198 (99%)	2 (1%)	0	100	100
1	F	200/230 (87%)	197 (98%)	3 (2%)	0	100	100
1	G	226/230 (98%)	215 (95%)	10 (4%)	1 (0%)	30	22
1	H	205/230 (89%)	201 (98%)	4 (2%)	0	100	100
All	All	1653/1840 (90%)	1619 (98%)	32 (2%)	2 (0%)	48	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	176	THR
1	G	188	THR

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	167/191 (87%)	162 (97%)	5 (3%)	36	30
1	B	167/191 (87%)	166 (99%)	1 (1%)	78	81
1	C	172/191 (90%)	166 (96%)	6 (4%)	32	24
1	D	177/191 (93%)	170 (96%)	7 (4%)	28	20
1	E	168/191 (88%)	165 (98%)	3 (2%)	51	50
1	F	167/191 (87%)	157 (94%)	10 (6%)	17	9
1	G	188/191 (98%)	175 (93%)	13 (7%)	14	7
1	H	171/191 (90%)	161 (94%)	10 (6%)	18	10
All	All	1377/1528 (90%)	1322 (96%)	55 (4%)	28	20

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	61	GLU
1	A	160	LYS
1	A	199	LEU
1	A	220	ARG
1	B	137	GLN
1	C	37	SER
1	C	66	GLN
1	C	87	GLN
1	C	133	GLN
1	C	199	LEU
1	C	207	ASN
1	D	8	SER
1	D	44	SER
1	D	87	GLN
1	D	165	LYS
1	D	207	ASN
1	D	220	ARG
1	D	221	GLU
1	E	56	SER
1	E	133	GLN
1	E	152	ILE
1	F	33	GLU
1	F	61	GLU
1	F	67	LEU
1	F	72	SER
1	F	87	GLN
1	F	162	LYS
1	F	167	ILE
1	F	207	ASN
1	F	216	ILE
1	F	220	ARG
1	G	33	GLU
1	G	34	SER
1	G	62	GLN
1	G	68	ASP
1	G	87	GLN
1	G	152	ILE
1	G	167	ILE
1	G	176	THR
1	G	177	GLU
1	G	184	LYS
1	G	199	LEU

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Mol	Chain	Res	Type
1	G	210	ASN
1	G	220	ARG
1	H	2	ILE
1	H	33	GLU
1	H	35	LYS
1	H	61	GLU
1	H	68	ASP
1	H	87	GLN
1	H	133	GLN
1	H	167	ILE
1	H	199	LEU
1	H	216	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	89	GLN
1	A	133	GLN
1	A	207	ASN
1	B	62	GLN
1	B	137	GLN
1	C	87	GLN
1	C	133	GLN
1	D	45	ASN
1	D	57	HIS
1	D	87	GLN
1	E	132	GLN
1	E	137	GLN
1	F	57	HIS
1	F	87	GLN
1	F	89	GLN
1	F	93	GLN
1	F	137	GLN
1	F	202	HIS
1	G	57	HIS
1	G	66	GLN
1	G	95	GLN
1	H	87	GLN
1	H	133	GLN

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	204/230 (88%)	0.05	3 (1%) 72 75	16, 27, 40, 50	0
1	B	204/230 (88%)	0.11	8 (3%) 43 46	15, 27, 42, 49	0
1	C	212/230 (92%)	0.29	9 (4%) 40 43	17, 29, 45, 60	0
1	D	220/230 (95%)	0.41	12 (5%) 30 32	17, 32, 49, 57	0
1	E	204/230 (88%)	0.25	5 (2%) 58 62	21, 32, 44, 56	0
1	F	204/230 (88%)	0.27	4 (1%) 65 69	20, 31, 43, 49	0
1	G	228/230 (99%)	0.83	28 (12%) 8 9	22, 39, 55, 61	0
1	H	209/230 (90%)	0.42	10 (4%) 35 38	21, 34, 45, 62	0
All	All	1685/1840 (91%)	0.34	79 (4%) 36 39	15, 31, 49, 62	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	179	TRP	6.2
1	E	191	PHE	5.0
1	H	178	PHE	5.0
1	C	179	TRP	4.5
1	H	206	ALA	4.4
1	B	191	PHE	4.0
1	D	176	THR	3.8
1	G	178	PHE	3.7
1	C	208	ILE	3.6
1	G	188	THR	3.5
1	H	207	ASN	3.4
1	G	176	THR	3.3
1	G	222	GLY	3.3
1	H	179	TRP	3.2
1	G	206	ALA	3.2
1	C	178	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	G	40	THR	3.1
1	F	172	GLY	3.1
1	G	43	LEU	3.1
1	G	121	VAL	3.1
1	D	211	GLY	3.1
1	D	186	LEU	3.0
1	H	211	GLY	3.0
1	D	181	THR	3.0
1	B	172	GLY	2.8
1	G	60	VAL	2.8
1	C	190	SER	2.8
1	G	189	SER	2.8
1	D	191	PHE	2.8
1	D	208	ILE	2.8
1	C	189	SER	2.7
1	A	207	ASN	2.7
1	H	177	GLU	2.7
1	D	38	THR	2.7
1	D	49	TYR	2.7
1	D	222	GLY	2.7
1	G	224	HIS	2.7
1	B	203	GLY	2.6
1	C	211	GLY	2.6
1	D	179	TRP	2.6
1	B	194	ALA	2.6
1	E	31	ARG	2.5
1	E	196	ASP	2.5
1	F	191	PHE	2.5
1	G	44	SER	2.5
1	D	206	ALA	2.5
1	E	172	GLY	2.4
1	G	191	PHE	2.4
1	H	161	GLY	2.4
1	B	193	SER	2.4
1	H	176	THR	2.4
1	C	205	LEU	2.3
1	E	61	GLU	2.3
1	D	175	ALA	2.3
1	G	207	ASN	2.3
1	C	207	ASN	2.3
1	G	210	ASN	2.3
1	A	196	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	68	ASP	2.3
1	G	63	LEU	2.3
1	B	210	ASN	2.3
1	C	21	ALA	2.3
1	G	18	LEU	2.3
1	G	177	GLU	2.2
1	G	225	HIS	2.2
1	G	226	HIS	2.2
1	G	1	LEU	2.2
1	A	191	PHE	2.1
1	G	228	HIS	2.1
1	H	133	GLN	2.1
1	G	49	TYR	2.1
1	B	196	ASP	2.1
1	B	44	SER	2.1
1	F	161	GLY	2.1
1	F	173	GLY	2.1
1	G	227	HIS	2.1
1	G	175	ALA	2.1
1	G	67	LEU	2.0
1	G	190	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.