



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 04:03 am BST

PDB ID : 3NKF
Title : Crystal structure of human ligand-free mature caspase-6 with intersubunit linker attached
Authors : Vaidya, S.; Hardy, J.A.
Deposited on : 2010-06-18
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

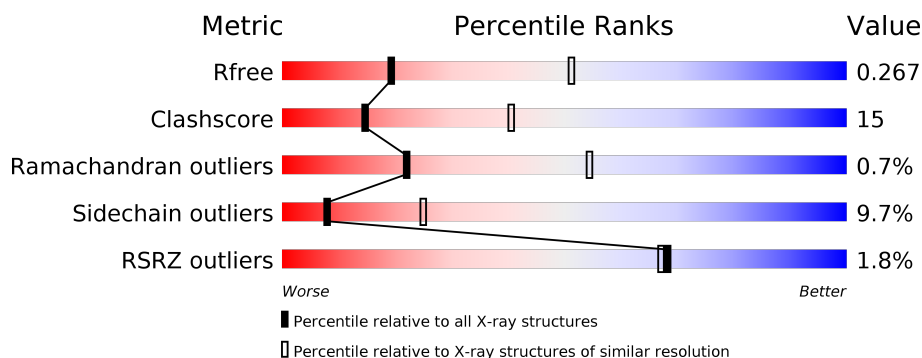
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on 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	277	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>18%</div> <div>•</div> <div>21%</div> </div> </div>
1	B	277	<div> <div>2%</div> <div> <div></div> <div>54%</div> <div>20%</div> <div>•</div> <div>23%</div> </div> </div>
1	C	277	<div> <div>%</div> <div> <div></div> <div>53%</div> <div>22%</div> <div>5%</div> <div>20%</div> </div> </div>
1	D	277	<div> <div>%</div> <div> <div></div> <div>55%</div> <div>21%</div> <div>•</div> <div>21%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6692 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 Caspase-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1665	1073	280	299	13			
1	B	213	Total	C	N	O	S	0	0	0
			1599	1019	273	294	13			
1	C	222	Total	C	N	O	S	0	0	0
			1722	1104	297	308	13			
1	D	220	Total	C	N	O	S	0	0	0
			1679	1075	286	305	13			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	EXPRESSION TAG	UNP P55212
A	294	HIS	-	EXPRESSION TAG	UNP P55212
A	295	HIS	-	EXPRESSION TAG	UNP P55212
A	296	HIS	-	EXPRESSION TAG	UNP P55212
A	297	HIS	-	EXPRESSION TAG	UNP P55212
A	298	HIS	-	EXPRESSION TAG	UNP P55212
A	299	HIS	-	EXPRESSION TAG	UNP P55212
B	23	MET	-	EXPRESSION TAG	UNP P55212
B	294	HIS	-	EXPRESSION TAG	UNP P55212
B	295	HIS	-	EXPRESSION TAG	UNP P55212
B	296	HIS	-	EXPRESSION TAG	UNP P55212
B	297	HIS	-	EXPRESSION TAG	UNP P55212
B	298	HIS	-	EXPRESSION TAG	UNP P55212
B	299	HIS	-	EXPRESSION TAG	UNP P55212
C	23	MET	-	EXPRESSION TAG	UNP P55212
C	294	HIS	-	EXPRESSION TAG	UNP P55212
C	295	HIS	-	EXPRESSION TAG	UNP P55212
C	296	HIS	-	EXPRESSION TAG	UNP P55212
C	297	HIS	-	EXPRESSION TAG	UNP P55212
C	298	HIS	-	EXPRESSION TAG	UNP P55212
C	299	HIS	-	EXPRESSION TAG	UNP P55212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	23	MET	-	EXPRESSION TAG	UNP P55212
D	294	HIS	-	EXPRESSION TAG	UNP P55212
D	295	HIS	-	EXPRESSION TAG	UNP P55212
D	296	HIS	-	EXPRESSION TAG	UNP P55212
D	297	HIS	-	EXPRESSION TAG	UNP P55212
D	298	HIS	-	EXPRESSION TAG	UNP P55212
D	299	HIS	-	EXPRESSION TAG	UNP P55212

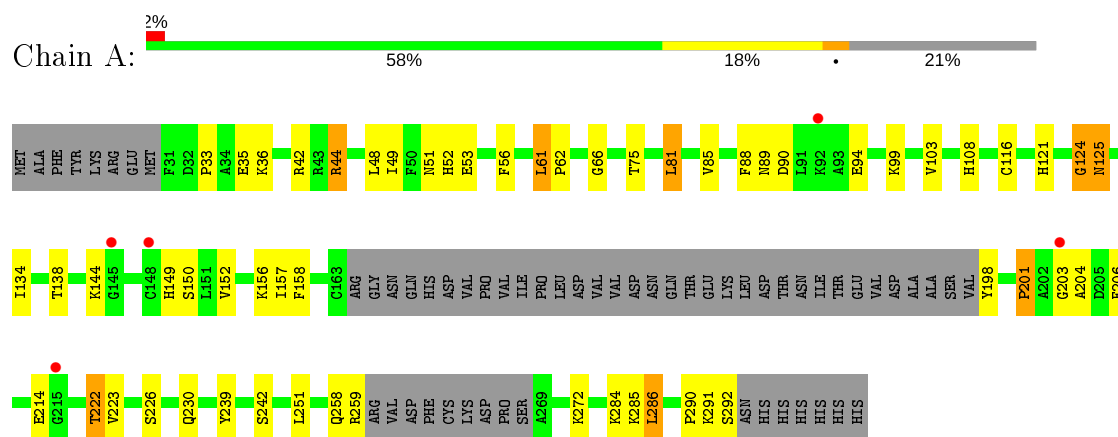
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total O 8 8	0	0
2	B	6	Total O 6 6	0	0
2	C	5	Total O 5 5	0	0
2	D	8	Total O 8 8	0	0

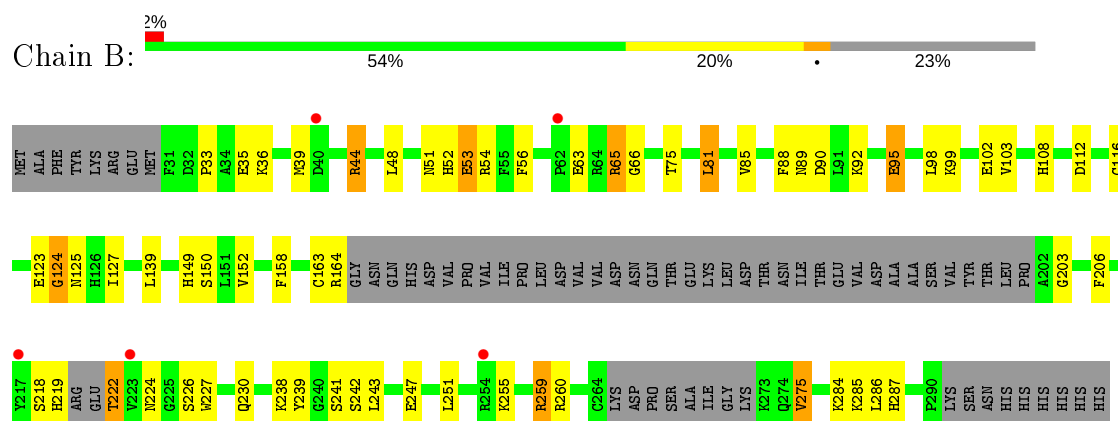
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

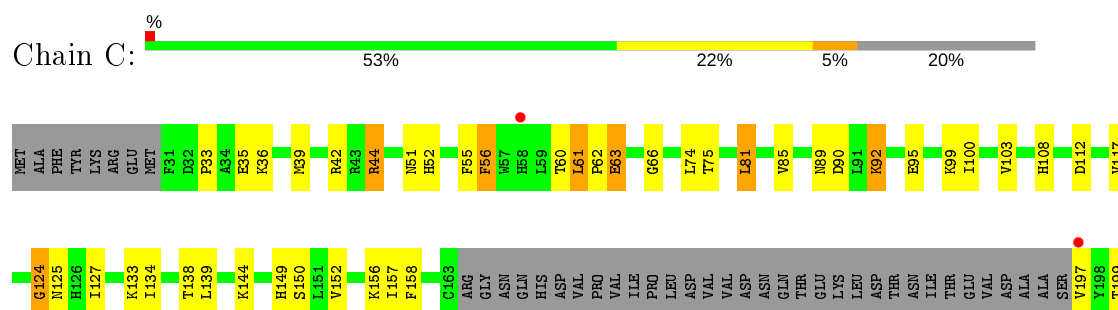
• Molecule 1: Caspase-6

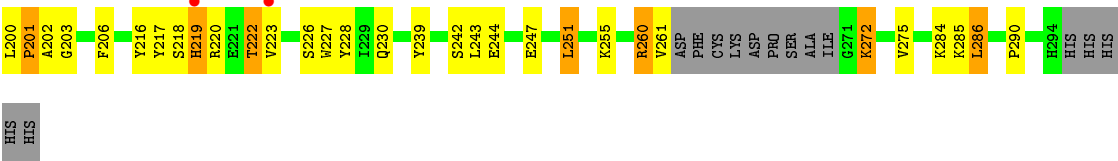


• Molecule 1: Caspase-6

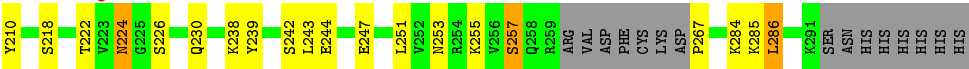


• Molecule 1: Caspase-6





● Molecule 1: Caspase-6



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.29Å 90.81Å 85.91Å 90.00° 91.04° 90.00°	Depositor
Resolution (Å)	29.50 – 2.90 29.52 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.50-2.90) 94.6 (29.52-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.216 , 0.270 0.214 , 0.267	Depositor DCC
R_{free} test set	1423 reflections (5.93%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 77.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.205 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6692	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.71% 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	0.51	0/1702	0.61	1/2298 (0.0%)
1	B	0.55	0/1631	0.61	0/2203
1	C	0.54	0/1760	0.64	0/2371
1	D	0.53	0/1715	0.63	1/2316 (0.0%)
All	All	0.53	0/6808	0.63	2/9188 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	267	PRO	N-CA-CB	6.09	110.61	103.30
1	A	201	PRO	N-CA-CB	5.83	110.30	103.30

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	1665	0	1559	38	0
1	B	1599	0	1465	63	0
1	C	1722	0	1638	61	0
1	D	1679	0	1584	46	0
2	A	8	0	0	2	0
2	B	6	0	0	2	0
2	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	8	0	0	0	0
All	All	6692	0	6246	197	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 15.

All (197) 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:287:HIS:HE1	2:B:9:HOH:O	1.49	0.93
1:C:219:HIS:HD2	1:C:227:TRP:HB2	1.34	0.92
1:B:287:HIS:CE1	2:B:9:HOH:O	2.24	0.90
1:C:52:HIS:HD2	1:C:90:ASP:OD2	1.53	0.90
1:C:219:HIS:CD2	1:C:227:TRP:HB2	2.12	0.84
1:A:52:HIS:HD2	1:A:90:ASP:OD2	1.61	0.83
1:B:227:TRP:CZ3	1:B:260:ARG:HA	2.14	0.82
1:C:200:LEU:N	1:C:201:PRO:HD2	1.95	0.82
1:C:203:GLY:HA3	1:C:206:PHE:CD1	2.14	0.81
1:D:52:HIS:HD2	1:D:90:ASP:OD2	1.62	0.81
1:C:133:LYS:HB3	1:C:197:VAL:CG2	2.11	0.81
1:C:44:ARG:HD2	1:C:81:LEU:O	1.82	0.79
1:B:203:GLY:HA3	1:B:206:PHE:CD1	2.17	0.78
1:A:204:ALA:HB2	1:B:275:VAL:HG21	1.68	0.75
1:C:239:TYR:CZ	1:D:33:PRO:HG3	2.21	0.75
1:B:92:LYS:CB	1:B:95:GLU:HG2	2.17	0.74
1:A:44:ARG:HD2	1:A:81:LEU:O	1.87	0.74
1:C:133:LYS:HG2	1:C:197:VAL:HG22	1.70	0.73
1:C:125:ASN:OD1	1:C:127:ILE:N	2.23	0.72
1:A:203:GLY:HA3	1:A:206:PHE:CD1	2.25	0.72
1:A:239:TYR:CZ	1:B:33:PRO:HG3	2.26	0.71
1:B:44:ARG:HD2	1:B:81:LEU:O	1.91	0.71
1:C:133:LYS:CG	1:C:197:VAL:HG22	2.23	0.69
1:B:92:LYS:O	1:B:95:GLU:HG3	1.92	0.69
1:D:44:ARG:HD2	1:D:81:LEU:O	1.93	0.68
1:A:291:LYS:HG3	2:A:11:HOH:O	1.93	0.67
1:C:133:LYS:CB	1:C:197:VAL:CG2	2.73	0.67
1:C:200:LEU:N	1:C:201:PRO:CD	2.58	0.67
1:B:125:ASN:OD1	1:B:127:ILE:N	2.28	0.66
1:B:75:THR:HG23	1:B:85:VAL:HG11	1.77	0.65
1:C:133:LYS:HB3	1:C:197:VAL:HG21	1.77	0.65
1:C:92:LYS:HD3	1:C:95:GLU:HB2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:PRO:HG3	1:D:239:TYR:CZ	2.32	0.65
1:B:158:PHE:HE1	1:B:206:PHE:HE2	1.44	0.64
1:B:92:LYS:CB	1:B:95:GLU:CG	2.77	0.63
1:C:217:TYR:CD2	1:C:272:LYS:HE3	2.34	0.63
1:A:56:PHE:HE1	1:A:124:GLY:H	1.44	0.63
1:C:51:ASN:HB3	1:C:89:ASN:HB3	1.79	0.63
1:B:158:PHE:HE1	1:B:206:PHE:CE2	2.18	0.62
1:B:218:SER:O	1:B:219:HIS:CG	2.52	0.62
1:B:227:TRP:CD1	1:B:259:ARG:NH1	2.69	0.61
1:A:125:ASN:ND2	1:A:125:ASN:H	1.98	0.60
1:B:158:PHE:CE1	1:B:206:PHE:HE2	2.19	0.60
1:C:52:HIS:CD2	1:C:90:ASP:OD2	2.45	0.60
1:D:200:LEU:N	1:D:201:PRO:HD3	2.16	0.60
1:A:290:PRO:HA	2:A:11:HOH:O	2.02	0.60
1:D:55:PHE:O	1:D:125:ASN:ND2	2.35	0.59
1:D:52:HIS:CD2	1:D:90:ASP:OD2	2.52	0.59
1:A:198:TYR:HA	1:A:214:GLU:CD	2.22	0.59
1:C:60:THR:OG1	1:C:63:GLU:HG2	2.02	0.59
1:B:36:LYS:HE3	1:B:287:HIS:HB2	1.85	0.58
1:D:60:THR:OG1	1:D:63:GLU:HG2	2.03	0.58
1:B:239:TYR:HB3	1:B:243:LEU:HG	1.85	0.58
1:C:100:ILE:HD11	1:C:139:LEU:CD2	2.34	0.58
1:C:219:HIS:HD2	1:C:227:TRP:CB	2.11	0.58
1:B:227:TRP:CH2	1:B:260:ARG:HA	2.39	0.57
1:B:238:LYS:HE3	1:C:290:PRO:O	2.03	0.57
1:B:227:TRP:CZ3	1:B:260:ARG:CA	2.87	0.57
1:C:55:PHE:O	1:C:125:ASN:ND2	2.37	0.57
1:B:52:HIS:HD2	1:B:90:ASP:OD2	1.87	0.57
1:C:133:LYS:CB	1:C:197:VAL:HG22	2.33	0.57
1:D:51:ASN:HB3	1:D:89:ASN:HB3	1.87	0.56
1:B:56:PHE:HE1	1:B:124:GLY:H	1.54	0.56
1:B:247:GLU:HA	1:B:247:GLU:OE1	2.06	0.56
1:C:133:LYS:HB3	1:C:197:VAL:HG22	1.85	0.56
1:A:134:ILE:O	1:A:138:THR:HG23	2.06	0.55
1:A:290:PRO:O	1:D:238:LYS:HE3	2.07	0.55
1:A:158:PHE:CE1	1:A:206:PHE:HE2	2.25	0.55
1:C:36:LYS:HG2	1:C:285:LYS:HB2	1.88	0.55
1:C:35:GLU:HG2	1:C:284:LYS:HG2	1.89	0.55
1:A:53:GLU:CD	1:A:121:HIS:HE2	2.10	0.54
1:C:260:ARG:HG2	1:C:261:VAL:H	1.72	0.54
1:C:99:LYS:O	1:C:103:VAL:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:226:SER:O	1:C:230:GLN:HG2	2.07	0.53
1:D:253:ASN:O	1:D:257:SER:HB3	2.08	0.53
1:B:36:LYS:HG2	1:B:285:LYS:HB2	1.90	0.53
1:C:199:THR:C	1:C:201:PRO:HD2	2.29	0.53
1:A:35:GLU:HG2	1:A:284:LYS:HG2	1.91	0.52
1:C:75:THR:HG23	1:C:85:VAL:HG11	1.91	0.52
1:A:52:HIS:CD2	1:A:90:ASP:OD2	2.52	0.52
1:D:61:LEU:CD1	1:D:65:ARG:HH11	2.22	0.52
1:A:75:THR:HG23	1:A:85:VAL:HG11	1.92	0.52
1:B:39:MET:SD	1:B:112:ASP:HB3	2.50	0.52
1:A:158:PHE:HE1	1:A:206:PHE:HE2	1.58	0.52
1:C:219:HIS:CD2	1:C:228:TYR:H	2.28	0.52
1:C:244:GLU:O	1:C:247:GLU:HB2	2.10	0.52
1:D:39:MET:SD	1:D:112:ASP:HB3	2.49	0.52
1:B:48:LEU:HD11	1:B:88:PHE:CZ	2.45	0.51
1:B:149:HIS:HA	1:B:152:VAL:HG23	1.92	0.51
1:D:239:TYR:HB3	1:D:243:LEU:HG	1.93	0.51
1:C:56:PHE:HE1	1:C:124:GLY:H	1.58	0.51
1:B:226:SER:O	1:B:230:GLN:HG2	2.10	0.51
1:A:157:ILE:HD11	1:A:286:LEU:HD11	1.91	0.50
1:B:125:ASN:C	1:B:125:ASN:OD1	2.50	0.50
1:B:116:CYS:HB3	1:B:158:PHE:CD2	2.47	0.50
1:D:146:ASP:OD1	1:D:146:ASP:N	2.45	0.50
1:C:200:LEU:O	1:C:202:ALA:N	2.45	0.50
1:D:61:LEU:HD12	1:D:65:ARG:HH11	1.76	0.50
1:A:108:HIS:H	1:A:150:SER:HB3	1.77	0.50
1:C:216:TYR:OH	1:D:203:GLY:HA2	2.12	0.50
1:B:51:ASN:HB3	1:B:89:ASN:HB3	1.94	0.50
1:B:203:GLY:HA3	1:B:206:PHE:HD1	1.74	0.49
1:B:203:GLY:HA3	1:B:206:PHE:CE1	2.46	0.49
1:B:51:ASN:ND2	1:B:53:GLU:OE2	2.44	0.49
1:D:226:SER:O	1:D:230:GLN:HG2	2.12	0.49
1:A:99:LYS:O	1:A:103:VAL:HG23	2.12	0.49
1:B:66:GLY:HA3	1:B:222:THR:O	2.12	0.49
1:A:51:ASN:HB3	1:A:89:ASN:HB3	1.93	0.49
1:D:75:THR:HG23	1:D:85:VAL:HG11	1.94	0.49
1:B:51:ASN:CG	1:B:53:GLU:OE2	2.51	0.49
1:D:203:GLY:HA3	1:D:206:PHE:CD1	2.47	0.48
1:A:226:SER:O	1:A:230:GLN:HG2	2.13	0.48
1:B:36:LYS:CE	1:B:287:HIS:HB2	2.43	0.48
1:C:125:ASN:OD1	1:C:125:ASN:C	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:TRP:HZ3	1:B:260:ARG:CB	2.27	0.48
1:C:133:LYS:HA	1:C:197:VAL:HG23	1.95	0.47
1:A:33:PRO:HG3	1:B:239:TYR:CZ	2.49	0.47
1:A:203:GLY:HA3	1:A:206:PHE:HD1	1.73	0.47
1:D:108:HIS:H	1:D:150:SER:HB3	1.78	0.47
1:B:238:LYS:HE2	1:B:239:TYR:OH	2.13	0.47
1:A:149:HIS:HA	1:A:152:VAL:HG23	1.97	0.47
1:B:108:HIS:H	1:B:150:SER:HB3	1.80	0.47
1:B:203:GLY:CA	1:B:206:PHE:CD1	2.94	0.47
1:D:56:PHE:HE1	1:D:124:GLY:H	1.62	0.47
1:D:104:SER:HB3	1:D:142:LEU:HD13	1.97	0.46
1:D:125:ASN:C	1:D:125:ASN:OD1	2.53	0.46
1:C:108:HIS:H	1:C:150:SER:HB3	1.81	0.46
1:B:203:GLY:O	1:B:206:PHE:HB2	2.16	0.46
1:D:36:LYS:HG2	1:D:285:LYS:HB2	1.98	0.46
1:B:98:LEU:O	1:B:102:GLU:HG3	2.16	0.46
1:C:149:HIS:HA	1:C:152:VAL:HG23	1.98	0.46
1:C:239:TYR:HB3	1:C:243:LEU:HG	1.97	0.45
1:B:35:GLU:HG2	1:B:284:LYS:HG2	1.98	0.45
1:D:108:HIS:H	1:D:150:SER:CB	2.30	0.45
1:A:66:GLY:HA3	1:A:222:THR:O	2.17	0.45
1:D:134:ILE:O	1:D:138:THR:HG23	2.16	0.45
1:A:158:PHE:HE1	1:A:206:PHE:CE2	2.33	0.45
1:C:39:MET:SD	1:C:112:ASP:HB3	2.57	0.45
1:D:224:ASN:O	1:D:230:GLN:OE1	2.35	0.45
1:C:144:LYS:HE2	1:C:144:LYS:HB2	1.69	0.44
1:B:56:PHE:CZ	1:B:123:GLU:HB2	2.52	0.44
1:D:158:PHE:CE1	1:D:206:PHE:HE2	2.36	0.44
1:B:65:ARG:HA	1:B:65:ARG:HD2	1.71	0.44
1:B:36:LYS:HB3	1:B:36:LYS:HE2	1.67	0.44
1:C:203:GLY:HA3	1:C:206:PHE:HD1	1.78	0.44
1:B:56:PHE:HZ	1:B:123:GLU:HB2	1.82	0.44
1:C:251:LEU:HD12	1:C:251:LEU:HA	1.79	0.44
1:B:99:LYS:O	1:B:103:VAL:HG23	2.18	0.44
1:D:158:PHE:HE1	1:D:206:PHE:CE2	2.36	0.44
1:A:116:CYS:HB3	1:A:158:PHE:CD2	2.53	0.43
1:A:144:LYS:HB2	1:A:144:LYS:HE2	1.68	0.43
1:A:203:GLY:O	1:A:206:PHE:HB2	2.19	0.43
1:C:100:ILE:HD11	1:C:139:LEU:HD21	2.01	0.43
1:C:108:HIS:H	1:C:150:SER:CB	2.32	0.43
1:C:66:GLY:HA3	1:C:222:THR:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:HIS:H	1:B:150:SER:CB	2.31	0.43
1:A:36:LYS:CG	1:A:285:LYS:HB2	2.49	0.42
1:B:92:LYS:CB	1:B:95:GLU:HG3	2.49	0.42
1:C:92:LYS:H	1:C:92:LYS:HG3	1.44	0.42
1:D:199:THR:HG23	1:D:210:TYR:CZ	2.55	0.42
1:C:61:LEU:HD22	1:C:61:LEU:HA	1.86	0.42
1:D:61:LEU:HA	1:D:61:LEU:HD22	1.89	0.42
1:C:61:LEU:N	1:C:62:PRO:CD	2.83	0.42
1:D:139:LEU:HD23	1:D:139:LEU:HA	1.81	0.42
1:D:99:LYS:O	1:D:103:VAL:HG23	2.20	0.42
1:A:108:HIS:H	1:A:150:SER:CB	2.32	0.42
1:B:163:CYS:O	1:B:164:ARG:O	2.37	0.42
1:C:239:TYR:CE1	1:D:33:PRO:HG3	2.54	0.42
1:A:48:LEU:HD12	1:A:49:ILE:N	2.34	0.42
1:B:218:SER:O	1:B:219:HIS:CD2	2.72	0.42
1:A:258:GLN:C	1:A:259:ARG:O	2.57	0.42
1:C:158:PHE:CE1	1:C:206:PHE:HE2	2.37	0.42
1:B:139:LEU:HD23	1:B:139:LEU:HA	1.87	0.42
1:D:61:LEU:N	1:D:62:PRO:CD	2.82	0.42
1:D:149:HIS:HA	1:D:152:VAL:HG23	2.02	0.41
1:D:157:ILE:HD11	1:D:286:LEU:HD11	2.02	0.41
1:B:203:GLY:CA	1:B:206:PHE:HD1	2.32	0.41
1:D:244:GLU:O	1:D:247:GLU:HB2	2.20	0.41
1:B:158:PHE:CE1	1:B:206:PHE:CE2	3.00	0.41
1:D:50:PHE:CD1	1:D:50:PHE:C	2.94	0.41
1:B:36:LYS:NZ	1:B:287:HIS:HB2	2.36	0.41
1:A:61:LEU:N	1:A:62:PRO:HD2	2.36	0.41
1:C:134:ILE:O	1:C:138:THR:HG23	2.19	0.41
1:D:61:LEU:O	1:D:65:ARG:HB2	2.21	0.41
1:B:203:GLY:C	1:B:206:PHE:HD1	2.24	0.41
1:D:32:ASP:HA	1:D:33:PRO:HD3	1.90	0.41
1:B:54:ARG:HD3	1:B:90:ASP:HB2	2.03	0.41
1:D:52:HIS:HE1	1:D:128:TYR:O	2.03	0.41
1:A:48:LEU:HD11	1:A:88:PHE:CZ	2.56	0.41
1:C:275:VAL:HG21	1:D:204:ALA:HB2	2.02	0.41
1:C:157:ILE:HD11	1:C:286:LEU:HD11	2.03	0.40
1:C:239:TYR:OH	1:D:33:PRO:HG3	2.20	0.40
1:B:227:TRP:CZ3	1:B:260:ARG:CB	3.03	0.40
1:C:74:LEU:HD13	1:C:117:VAL:HG11	2.03	0.40
1:A:158:PHE:CE1	1:A:206:PHE:CE2	3.08	0.40
1:D:35:GLU:HG2	1:D:284:LYS:HG2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:PHE:CD2	1:C:127:ILE:CG2	3.05	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	213/277 (77%)	193 (91%)	18 (8%)	2 (1%)	17	48
1	B	205/277 (74%)	193 (94%)	11 (5%)	1 (0%)	29	61
1	C	216/277 (78%)	199 (92%)	15 (7%)	2 (1%)	17	48
1	D	214/277 (77%)	197 (92%)	16 (8%)	1 (0%)	29	61
All	All	848/1108 (76%)	782 (92%)	60 (7%)	6 (1%)	22	54

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	201	PRO
1	A	124	GLY
1	B	124	GLY
1	C	124	GLY
1	D	124	GLY
1	A	201	PRO

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	163/245 (66%)	149 (91%)	14 (9%)	10	30
1	B	156/245 (64%)	141 (90%)	15 (10%)	8	25
1	C	173/245 (71%)	154 (89%)	19 (11%)	6	19
1	D	169/245 (69%)	153 (90%)	16 (10%)	8	26
All	All	661/980 (67%)	597 (90%)	64 (10%)	8	25

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	44	ARG
1	A	61	LEU
1	A	81	LEU
1	A	94	GLU
1	A	125	ASN
1	A	156	LYS
1	A	222	THR
1	A	223	VAL
1	A	242	SER
1	A	251	LEU
1	A	272	LYS
1	A	286	LEU
1	A	292	SER
1	B	44	ARG
1	B	53	GLU
1	B	63	GLU
1	B	65	ARG
1	B	81	LEU
1	B	95	GLU
1	B	222	THR
1	B	224	ASN
1	B	241	SER
1	B	242	SER
1	B	251	LEU
1	B	255	LYS
1	B	259	ARG
1	B	275	VAL
1	B	286	LEU
1	C	42	ARG
1	C	44	ARG

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Mol	Chain	Res	Type
1	C	56	PHE
1	C	61	LEU
1	C	63	GLU
1	C	81	LEU
1	C	92	LYS
1	C	156	LYS
1	C	218	SER
1	C	219	HIS
1	C	220	ARG
1	C	222	THR
1	C	223	VAL
1	C	242	SER
1	C	251	LEU
1	C	255	LYS
1	C	260	ARG
1	C	272	LYS
1	C	286	LEU
1	D	44	ARG
1	D	61	LEU
1	D	81	LEU
1	D	98	LEU
1	D	103	VAL
1	D	146	ASP
1	D	156	LYS
1	D	200	LEU
1	D	218	SER
1	D	222	THR
1	D	224	ASN
1	D	242	SER
1	D	251	LEU
1	D	255	LYS
1	D	257	SER
1	D	286	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	125	ASN
1	A	230	GLN
1	A	258	GLN
1	B	52	HIS

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Mol	Chain	Res	Type
1	B	108	HIS
1	B	230	GLN
1	B	258	GLN
1	C	52	HIS
1	C	219	HIS
1	C	258	GLN
1	D	52	HIS
1	D	258	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 carbohydrates 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	219/277 (79%)	0.03	5 (2%) 60 58	57, 85, 143, 173	0
1	B	213/277 (76%)	0.03	5 (2%) 60 58	56, 89, 139, 181	0
1	C	222/277 (80%)	-0.16	4 (1%) 68 67	29, 69, 128, 214	0
1	D	220/277 (79%)	-0.18	2 (0%) 84 84	30, 71, 127, 191	0
All	All	874/1108 (78%)	-0.07	16 (1%) 68 67	29, 79, 138, 214	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	62	PRO	4.1
1	A	145	GLY	3.8
1	A	215	GLY	3.2
1	A	148	CYS	3.2
1	D	223	VAL	3.2
1	B	40	ASP	3.1
1	B	217	TYR	2.7
1	C	58	HIS	2.6
1	C	197	VAL	2.5
1	B	223	VAL	2.4
1	D	98	LEU	2.3
1	A	203	GLY	2.2
1	C	223	VAL	2.2
1	A	92	LYS	2.2
1	C	219	HIS	2.2
1	B	254	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.