



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:46 PM UTC

PDB ID : 3SOH / pdb_00003soh
Title : Architecture of the Flagellar Rotor
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Deposited on : 2011-06-30
Resolution : 3.50 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

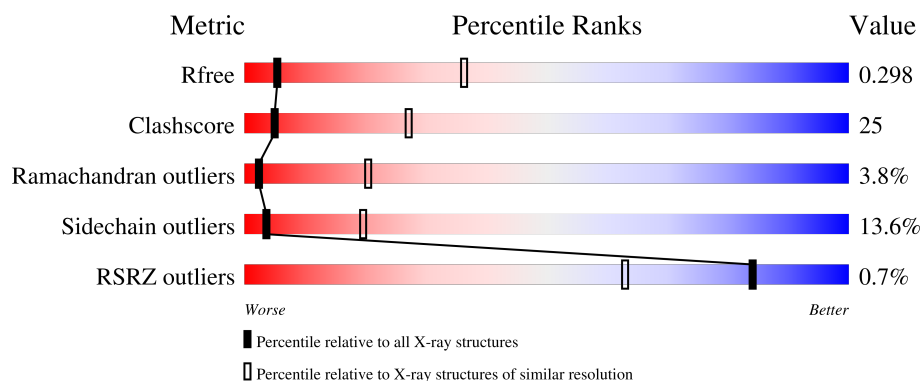
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	188	
1	C	188	
2	B	80	
2	D	80	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4285 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 Flagellar motor switch protein FliM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1535	1003	240	283	9			
1	C	188	Total	C	N	O	S	0	0	0
			1535	1003	240	283	9			

- Molecule 2 is a protein called Flagellar motor switch protein FliG.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	80	Total	C	N	O	0	0	0
			607	387	101	119			
2	D	80	Total	C	N	O	0	0	0
			607	387	101	119			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	113	GLY	-	cloning artifact	UNP Q9WY63
B	114	SER	-	cloning artifact	UNP Q9WY63
B	115	HIS	-	cloning artifact	UNP Q9WY63
B	116	MET	-	cloning artifact	UNP Q9WY63
B	?	-	VAL	deletion	UNP Q9WY63
D	113	GLY	-	cloning artifact	UNP Q9WY63
D	114	SER	-	cloning artifact	UNP Q9WY63
D	115	HIS	-	cloning artifact	UNP Q9WY63
D	116	MET	-	cloning artifact	UNP Q9WY63
D	?	-	VAL	deletion	UNP Q9WY63

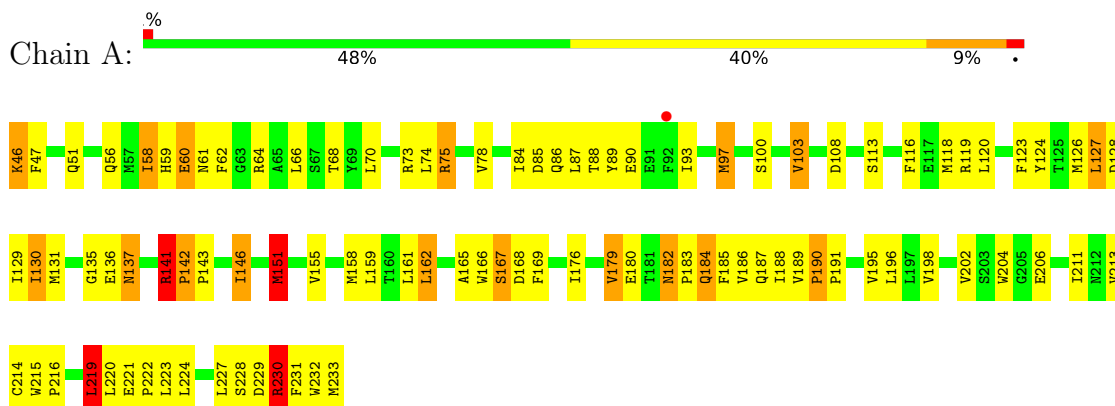
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	O	0	0
			1	1		

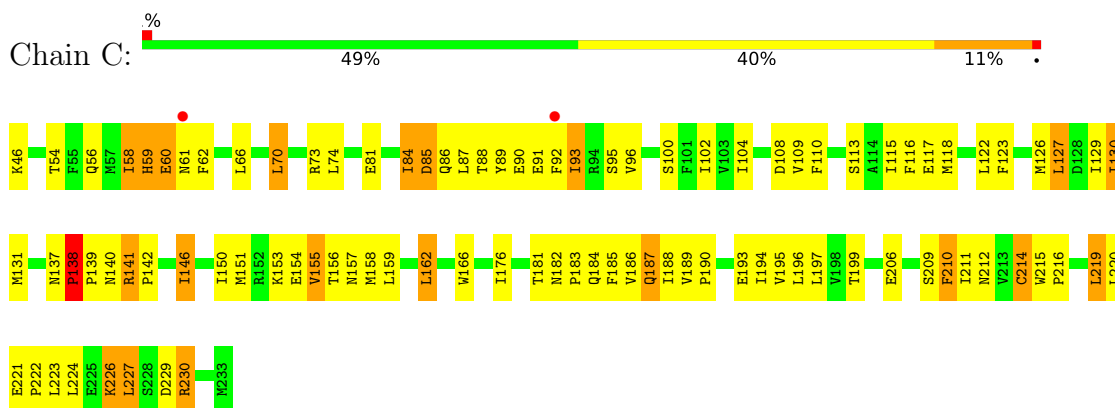
3 Residue-property plots

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.

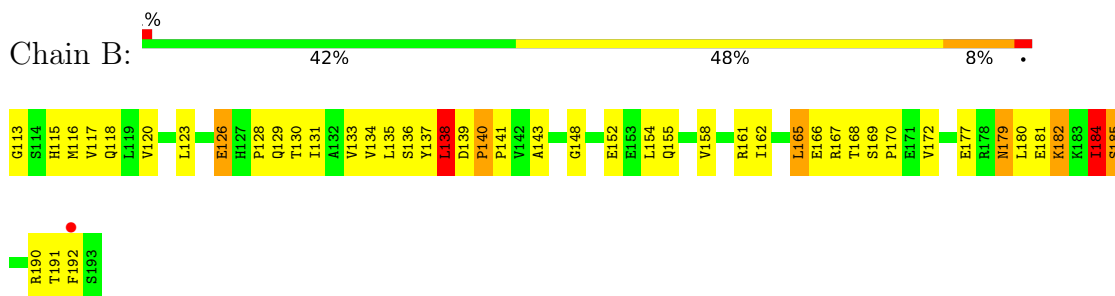
• Molecule 1: Flagellar motor switch protein FliM



• Molecule 1: Flagellar motor switch protein FliM



• Molecule 2: Flagellar motor switch protein FliG



• Molecule 2: Flagellar motor switch protein FliG

Chain D: 40% 46% 14%



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	91.39Å 91.39Å 226.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.21 – 3.50 26.21 – 3.50	Depositor EDS
% Data completeness (in resolution range)	87.2 (26.21-3.50) 86.9 (26.21-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 3.32Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.253 , 0.301 0.258 , 0.298	Depositor DCC
R_{free} test set	1608 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	88.3	Xtriage
Anisotropy	0.737	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 147.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4285	wwPDB-VP
Average B, all atoms (Å ²)	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% 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	0.66	3/1576 (0.2%)	1.20	17/2143 (0.8%)
1	C	0.55	0/1576	1.14	14/2143 (0.7%)
2	B	0.78	1/616 (0.2%)	1.27	4/838 (0.5%)
2	D	0.71	0/616	1.25	4/838 (0.5%)
All	All	0.65	4/4384 (0.1%)	1.20	39/5962 (0.7%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	229	ASP	C-O	7.51	1.33	1.24
2	B	165	LEU	N-CA	5.58	1.53	1.46
1	A	228	SER	C-N	5.50	1.41	1.33
1	A	151	MET	SD-CE	-5.19	1.66	1.79

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	138	PRO	CA-C-N	10.25	130.34	119.89
1	C	138	PRO	C-N-CA	10.25	130.34	119.89
1	A	230	ARG	N-CA-C	10.05	123.17	111.11
1	A	190	PRO	CA-C-N	9.04	131.14	119.84
1	A	190	PRO	C-N-CA	9.04	131.14	119.84
2	D	190	ARG	N-CA-C	7.91	120.81	111.71
2	D	140	PRO	N-CA-C	7.76	120.17	110.70
1	C	229	ASP	N-CA-C	7.36	121.56	108.24
1	C	140	ASN	N-CA-C	6.92	121.44	112.92
2	B	140	PRO	N-CA-C	6.77	118.96	110.70
1	C	206	GLU	N-CA-C	-6.50	105.65	113.97
1	A	142	PRO	CA-C-N	6.14	126.08	119.76
1	A	142	PRO	C-N-CA	6.14	126.08	119.76
1	C	190	PRO	CA-C-N	6.14	127.52	119.84
1	C	190	PRO	C-N-CA	6.14	127.52	119.84
2	B	184	ILE	N-CA-C	-6.10	100.84	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	185	SER	N-CA-C	6.01	123.61	110.80
1	A	219	LEU	N-CA-C	-5.98	106.96	114.56
1	A	162	LEU	N-CA-C	-5.96	104.69	111.07
1	A	182	ASN	CA-C-N	5.85	125.55	119.05
1	A	182	ASN	C-N-CA	5.85	125.55	119.05
1	C	193	GLU	N-CA-C	5.80	118.46	110.35
2	D	189	SER	N-CA-C	5.75	117.97	109.69
1	C	210	PHE	N-CA-C	5.72	118.53	109.50
1	A	141	ARG	CA-C-N	5.59	126.14	120.38
1	A	141	ARG	C-N-CA	5.59	126.14	120.38
1	A	47	PHE	N-CA-C	5.57	118.63	109.72
1	A	60	GLU	N-CA-C	-5.51	105.27	111.28
1	A	58	ILE	CB-CA-C	-5.46	104.98	111.81
1	C	197	LEU	N-CA-C	5.46	117.20	108.41
1	A	229	ASP	O-C-N	-5.38	115.44	122.59
1	A	229	ASP	CB-CA-C	5.34	121.04	110.42
1	C	156	THR	N-CA-C	-5.22	105.67	111.36
1	C	60	GLU	N-CA-C	-5.17	105.73	111.36
1	A	206	GLU	N-CA-C	-5.13	106.87	113.02
1	C	227	LEU	N-CA-C	5.13	117.48	108.56
2	D	126	GLU	N-CA-C	5.11	116.89	110.24
1	C	157	ASN	N-CA-C	-5.07	105.65	111.07
2	B	180	LEU	N-CA-C	-5.01	105.73	111.14

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	1535	0	1516	81	0
1	C	1535	0	1516	69	0
2	B	607	0	615	35	0
2	D	607	0	615	35	0
3	B	1	0	0	0	0
All	All	4285	0	4262	214	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 25.

All (214) 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:195:VAL:HG11	1:A:214:CYS:SG	2.01	1.01
1:A:113:SER:HB2	1:A:188:ILE:HD11	1.46	0.96
1:C:113:SER:HB2	1:C:188:ILE:HD11	1.57	0.87
2:D:134:VAL:HA	2:D:192:PHE:HZ	1.39	0.85
2:D:138:LEU:HD22	2:D:191:THR:HG21	1.58	0.85
1:C:59:HIS:CE1	1:C:166:TRP:HE1	1.96	0.83
1:A:227:LEU:HD22	2:D:170:PRO:HB3	1.62	0.81
1:C:146:ILE:HD13	1:C:146:ILE:H	1.47	0.77
1:A:59:HIS:HE1	1:A:215:TRP:HE1	1.31	0.77
1:A:116:PHE:HZ	1:A:158:MET:HE3	1.50	0.77
1:C:122:LEU:HD23	1:C:211:ILE:HD11	1.65	0.77
1:C:188:ILE:HG23	1:C:189:VAL:HG22	1.67	0.76
1:A:84:ILE:HD11	1:A:196:LEU:HD11	1.67	0.75
1:A:220:LEU:O	1:A:224:LEU:HB2	1.87	0.74
1:C:93:ILE:HD11	1:C:186:VAL:HG23	1.68	0.74
2:D:118:GLN:NE2	2:D:184:ILE:HA	2.05	0.72
1:A:89:TYR:CZ	1:A:188:ILE:HG22	2.25	0.71
1:C:220:LEU:O	1:C:224:LEU:HB2	1.91	0.70
2:B:167:ARG:O	2:B:168:THR:HG23	1.92	0.70
2:B:137:TYR:HD1	2:B:192:PHE:HE1	1.41	0.68
1:C:113:SER:HB2	1:C:188:ILE:CD1	2.23	0.68
1:C:122:LEU:CD2	1:C:211:ILE:HD11	2.22	0.68
1:C:199:THR:HG23	1:C:212:ASN:HD21	1.58	0.68
1:A:155:VAL:O	1:A:159:LEU:HB2	1.94	0.68
1:A:146:ILE:HD13	1:A:146:ILE:H	1.57	0.67
1:C:195:VAL:HG11	1:C:214:CYS:SG	2.35	0.67
2:B:138:LEU:HA	2:B:191:THR:HG21	1.76	0.67
1:A:116:PHE:CZ	1:A:158:MET:HE3	2.30	0.67
1:C:86:GLN:O	1:C:87:LEU:HD12	1.95	0.67
2:D:134:VAL:HA	2:D:192:PHE:CZ	2.26	0.67
1:C:84:ILE:HD11	1:C:196:LEU:HD11	1.77	0.66
1:A:186:VAL:HG23	1:A:187:GLN:H	1.61	0.66
1:C:196:LEU:HB3	1:C:215:TRP:HB2	1.79	0.65
2:B:128:PRO:HB2	2:B:161:ARG:HE	1.62	0.65
1:A:100:SER:HA	1:A:120:LEU:HG	1.79	0.64
1:C:66:LEU:HG	1:C:70:LEU:CD2	2.28	0.64
1:C:127:LEU:CD2	1:C:131:MET:SD	2.86	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:GLN:O	1:C:60:GLU:HB2	1.98	0.64
1:C:141:ARG:HB2	1:C:142:PRO:HD2	1.80	0.64
1:C:59:HIS:CE1	1:C:215:TRP:HE1	2.17	0.63
1:A:186:VAL:HG23	1:A:187:GLN:N	2.14	0.63
2:B:137:TYR:HD1	2:B:192:PHE:CE1	2.17	0.63
1:C:199:THR:HG23	1:C:212:ASN:ND2	2.13	0.62
2:D:129:GLN:HE21	2:D:161:ARG:HH11	1.46	0.62
2:B:131:ILE:HG22	2:B:135:LEU:HD12	1.82	0.62
1:C:155:VAL:O	1:C:159:LEU:HB2	2.01	0.61
2:D:122:PHE:HE2	2:D:180:LEU:CD1	2.15	0.60
1:A:56:GLN:O	1:A:60:GLU:HB2	2.02	0.60
2:B:140:PRO:HB2	2:B:141:PRO:HD3	1.83	0.60
1:C:89:TYR:CZ	1:C:188:ILE:HG22	2.36	0.60
1:C:104:ILE:HG22	1:C:115:ILE:HG12	1.84	0.60
1:C:166:TRP:HZ3	1:C:223:LEU:HD11	1.67	0.60
2:B:133:VAL:O	2:B:136:SER:HB2	2.01	0.60
2:D:141:PRO:O	2:D:145:GLN:HG3	2.02	0.60
1:A:128:ASP:OD1	1:A:135:GLY:HA2	2.01	0.60
1:C:127:LEU:HD22	1:C:131:MET:SD	2.42	0.60
1:A:59:HIS:CE1	1:A:215:TRP:HE1	2.15	0.59
1:C:96:VAL:HG11	1:C:102:ILE:HD11	1.84	0.59
1:A:123:PHE:CE1	1:A:151:MET:HG2	2.38	0.59
1:A:183:PRO:C	1:A:185:PHE:H	2.09	0.59
1:C:183:PRO:C	1:C:185:PHE:H	2.10	0.59
2:B:129:GLN:HE21	2:B:161:ARG:NH1	2.01	0.58
1:C:216:PRO:HG2	1:C:219:LEU:HB2	1.86	0.57
2:B:129:GLN:HE21	2:B:161:ARG:HH11	1.51	0.57
1:A:221:GLU:HG3	1:A:222:PRO:HD3	1.87	0.57
1:A:188:ILE:HG23	1:A:189:VAL:HG22	1.86	0.56
2:D:137:TYR:HB2	2:D:192:PHE:CE1	2.41	0.56
1:A:90:GLU:HG2	1:A:191:PRO:HB2	1.88	0.56
1:A:189:VAL:HB	1:A:190:PRO:HD2	1.87	0.56
1:A:216:PRO:HG2	1:A:219:LEU:HB2	1.87	0.56
2:D:181:GLU:HG3	2:D:187:PHE:CB	2.36	0.56
2:D:184:ILE:HG22	2:D:184:ILE:O	2.06	0.56
1:C:87:LEU:O	1:C:194:ILE:HD12	2.07	0.55
1:A:74:LEU:HD21	1:A:126:MET:HG2	1.87	0.55
1:A:223:LEU:HD23	1:A:223:LEU:H	1.72	0.55
1:A:59:HIS:HE1	1:A:215:TRP:NE1	2.04	0.54
1:A:90:GLU:HG2	1:A:191:PRO:CB	2.37	0.54
1:C:150:ILE:O	1:C:153:LYS:HG2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:VAL:CG2	1:A:179:VAL:HG13	2.36	0.54
2:B:134:VAL:HA	2:B:192:PHE:HZ	1.72	0.54
2:B:136:SER:OG	2:B:165:LEU:HD13	2.07	0.54
2:D:160:LYS:O	2:D:163:ALA:HB3	2.08	0.54
2:D:127:HIS:ND1	2:D:128:PRO:HD2	2.22	0.54
1:C:90:GLU:HG3	1:C:91:GLU:N	2.23	0.53
1:A:113:SER:HB2	1:A:188:ILE:CD1	2.30	0.53
1:A:62:PHE:O	1:A:66:LEU:HB2	2.08	0.53
2:D:152:GLU:CD	2:D:152:GLU:H	2.17	0.53
2:B:116:MET:O	2:B:120:VAL:HG23	2.09	0.52
2:B:148:GLY:HA2	2:B:155:GLN:NE2	2.24	0.52
2:B:170:PRO:HB3	1:C:227:LEU:HD22	1.92	0.52
1:C:116:PHE:HZ	1:C:158:MET:HE3	1.74	0.52
1:A:74:LEU:HD12	1:A:78:VAL:HG21	1.91	0.52
1:C:62:PHE:CD1	1:C:162:LEU:HA	2.45	0.52
1:A:182:ASN:HB3	1:A:185:PHE:CD2	2.44	0.52
1:C:92:PHE:O	1:C:96:VAL:HG23	2.09	0.52
2:B:169:SER:CB	2:B:172:VAL:HG12	2.41	0.51
1:C:58:ILE:HG21	1:C:166:TRP:CD1	2.45	0.51
1:A:128:ASP:CG	1:A:135:GLY:HA2	2.36	0.51
2:B:129:GLN:NE2	2:B:161:ARG:HH11	2.08	0.50
1:A:97:MET:O	1:A:100:SER:HB2	2.12	0.50
2:B:152:GLU:CD	2:B:152:GLU:H	2.19	0.50
2:D:115:HIS:O	2:D:119:LEU:HB2	2.11	0.50
2:B:169:SER:HB3	2:B:172:VAL:HG12	1.93	0.50
2:D:116:MET:O	2:D:120:VAL:HG23	2.12	0.50
1:A:166:TRP:HZ3	1:A:223:LEU:HD11	1.76	0.50
1:A:131:MET:HG2	2:B:172:VAL:HG21	1.94	0.50
1:A:120:LEU:HD22	1:A:124:TYR:CE1	2.47	0.49
2:B:169:SER:HB3	2:B:172:VAL:CG1	2.42	0.49
1:A:230:ARG:NH1	1:A:231:PHE:HD2	2.10	0.49
2:B:113:GLY:O	2:B:117:VAL:HG23	2.13	0.49
1:C:186:VAL:HG23	1:C:187:GLN:H	1.78	0.49
2:D:177:GLU:C	2:D:179:ASN:H	2.18	0.49
1:A:61:ASN:HA	1:A:64:ARG:HG3	1.93	0.49
1:A:93:ILE:HD11	1:A:186:VAL:CG2	2.43	0.49
1:A:84:ILE:HD11	1:A:196:LEU:CD1	2.40	0.49
2:D:122:PHE:HE2	2:D:180:LEU:HD12	1.78	0.49
1:C:56:GLN:HE21	1:C:84:ILE:HG22	1.78	0.48
2:B:169:SER:OG	2:B:172:VAL:HG12	2.13	0.48
1:C:118:MET:HG3	1:C:211:ILE:CD1	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:137:TYR:HB2	2:D:192:PHE:HE1	1.79	0.48
1:A:59:HIS:CE1	1:A:215:TRP:NE1	2.80	0.48
2:B:182:LYS:NZ	2:B:182:LYS:HB3	2.29	0.48
1:C:74:LEU:HD21	1:C:126:MET:HG2	1.96	0.48
2:B:158:VAL:O	2:B:162:ILE:HG13	2.14	0.48
1:A:131:MET:HE1	2:B:130:THR:HA	1.96	0.47
2:D:129:GLN:NE2	2:D:161:ARG:HD3	2.28	0.47
2:D:129:GLN:O	2:D:133:VAL:HG23	2.13	0.47
2:D:130:THR:O	2:D:134:VAL:HG23	2.14	0.47
1:A:141:ARG:HB2	1:A:142:PRO:HD2	1.96	0.47
1:A:120:LEU:HD22	1:A:124:TYR:HE1	1.80	0.47
2:D:140:PRO:HB2	2:D:141:PRO:HD3	1.96	0.47
2:B:128:PRO:HB2	2:B:161:ARG:NE	2.28	0.47
1:A:51:GLN:HE22	1:A:224:LEU:HD11	1.78	0.47
1:A:196:LEU:HB3	1:A:215:TRP:HB2	1.95	0.47
1:C:183:PRO:O	1:C:186:VAL:HG22	2.14	0.47
1:A:129:ILE:C	1:A:131:MET:H	2.22	0.46
2:D:137:TYR:HD1	2:D:192:PHE:HE1	1.63	0.46
1:A:190:PRO:HA	1:A:191:PRO:HD2	1.61	0.46
1:C:73:ARG:HE	1:C:73:ARG:HA	1.81	0.46
1:A:86:GLN:C	1:A:87:LEU:HD12	2.40	0.46
1:A:232:TRP:HE3	1:A:233:MET:HB2	1.81	0.46
1:A:232:TRP:CE3	1:A:233:MET:HB2	2.50	0.46
2:D:127:HIS:O	2:D:131:ILE:HG13	2.16	0.45
1:A:182:ASN:O	1:A:185:PHE:HB2	2.17	0.45
2:B:136:SER:OG	2:B:165:LEU:CD1	2.64	0.45
1:C:66:LEU:O	1:C:70:LEU:HD22	2.16	0.45
1:C:129:ILE:C	1:C:131:MET:H	2.23	0.45
2:D:177:GLU:C	2:D:179:ASN:N	2.73	0.45
1:C:58:ILE:O	1:C:61:ASN:N	2.45	0.45
1:C:74:LEU:HA	1:C:129:ILE:HG21	1.98	0.45
1:C:123:PHE:CZ	1:C:151:MET:HG3	2.50	0.45
1:C:182:ASN:O	1:C:185:PHE:HB2	2.16	0.45
1:C:221:GLU:HG3	1:C:222:PRO:HD3	1.98	0.45
1:A:123:PHE:C	1:A:123:PHE:CD2	2.95	0.45
1:C:151:MET:HA	1:C:154:GLU:HB2	1.99	0.45
2:D:129:GLN:HE21	2:D:161:ARG:HD3	1.81	0.45
1:A:58:ILE:HG21	1:A:166:TRP:CD1	2.53	0.44
1:C:146:ILE:HD13	1:C:146:ILE:N	2.24	0.44
1:A:74:LEU:O	1:A:75:ARG:HB2	2.16	0.44
2:B:140:PRO:O	2:B:143:ALA:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ILE:HG12	1:C:85:ASP:N	2.33	0.44
1:C:89:TYR:CZ	1:C:93:ILE:HD12	2.53	0.44
1:A:188:ILE:HG23	1:A:189:VAL:N	2.33	0.44
2:B:154:LEU:O	2:B:155:GLN:C	2.60	0.43
2:D:157:GLU:O	2:D:161:ARG:HG3	2.18	0.43
1:A:103:VAL:HG22	1:A:179:VAL:HG13	2.00	0.43
1:A:119:ARG:HG3	1:A:119:ARG:HH11	1.83	0.43
2:B:177:GLU:C	2:B:179:ASN:H	2.27	0.43
1:C:146:ILE:HD11	2:D:126:GLU:OE1	2.18	0.43
1:C:183:PRO:C	1:C:185:PHE:N	2.77	0.43
1:A:141:ARG:CB	1:A:142:PRO:HD2	2.48	0.43
1:A:161:LEU:H	1:A:161:LEU:HG	1.54	0.43
1:A:169:PHE:N	1:A:169:PHE:CD1	2.87	0.43
1:A:183:PRO:C	1:A:185:PHE:N	2.75	0.43
2:B:118:GLN:HB3	2:B:184:ILE:HG13	2.00	0.43
2:D:169:SER:OG	2:D:172:VAL:HG13	2.19	0.43
1:A:127:LEU:HD22	1:A:131:MET:SD	2.58	0.43
1:C:138:PRO:HA	1:C:139:PRO:HD2	1.72	0.43
1:A:58:ILE:HG23	1:A:165:ALA:O	2.18	0.42
1:A:221:GLU:HG3	1:A:222:PRO:CD	2.49	0.42
1:A:198:VAL:HB	1:A:213:VAL:HB	2.01	0.42
1:A:221:GLU:N	1:A:222:PRO:HD2	2.34	0.42
1:C:100:SER:O	1:C:181:THR:HA	2.19	0.42
1:C:102:ILE:HG12	1:C:117:GLU:HG3	2.00	0.42
1:C:183:PRO:O	1:C:185:PHE:N	2.52	0.42
1:C:222:PRO:O	1:C:226:LYS:HE2	2.20	0.42
1:A:103:VAL:HG23	1:A:179:VAL:HG13	2.00	0.42
1:C:129:ILE:O	1:C:131:MET:N	2.53	0.42
1:A:142:PRO:HA	1:A:143:PRO:HD2	1.82	0.42
1:A:230:ARG:HH12	1:A:231:PHE:HD2	1.66	0.42
1:C:54:THR:O	1:C:58:ILE:HG13	2.20	0.42
1:C:141:ARG:CB	1:C:142:PRO:HD2	2.46	0.42
1:A:118:MET:HG3	1:A:211:ILE:CD1	2.50	0.42
1:A:155:VAL:HA	1:A:158:MET:HE2	2.01	0.42
1:C:151:MET:O	1:C:155:VAL:HG13	2.19	0.42
1:A:129:ILE:O	1:A:131:MET:N	2.53	0.41
1:A:137:ASN:O	1:A:137:ASN:ND2	2.53	0.41
1:C:209:SER:OG	1:C:210:PHE:N	2.53	0.41
2:D:177:GLU:O	2:D:179:ASN:N	2.53	0.41
1:A:46:LYS:HD3	1:A:46:LYS:N	2.34	0.41
2:D:137:TYR:CD1	2:D:192:PHE:HE1	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:SER:O	2:D:171:GLU:HG3	2.21	0.41
1:C:59:HIS:HE1	1:C:166:TRP:HE1	1.62	0.41
1:A:126:MET:HE2	1:A:126:MET:HB3	1.93	0.41
1:A:90:GLU:CG	1:A:191:PRO:HB2	2.51	0.41
1:C:127:LEU:HA	1:C:130:ILE:HG12	2.02	0.41
2:B:123:LEU:HD23	2:B:123:LEU:HA	1.86	0.41
2:B:126:GLU:HB3	2:B:130:THR:HB	2.03	0.41
2:B:137:TYR:CD1	2:B:192:PHE:HE1	2.29	0.41
1:C:73:ARG:HA	1:C:73:ARG:NE	2.36	0.41
1:A:59:HIS:CE1	1:A:166:TRP:HE1	2.39	0.41
2:D:173:VAL:C	2:D:175:GLU:H	2.29	0.41
2:D:139:ASP:O	2:D:143:ALA:N	2.52	0.40
1:C:86:GLN:C	1:C:87:LEU:HD12	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	186/188 (99%)	164 (88%)	19 (10%)	3 (2%)	7	36
1	C	186/188 (99%)	165 (89%)	13 (7%)	8 (4%)	2	18
2	B	78/80 (98%)	57 (73%)	15 (19%)	6 (8%)	1	8
2	D	78/80 (98%)	63 (81%)	12 (15%)	3 (4%)	2	20
All	All	528/536 (98%)	449 (85%)	59 (11%)	20 (4%)	2	20

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	185	SER
1	A	130	ILE

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Mol	Chain	Res	Type
2	B	126	GLU
2	B	190	ARG
1	C	184	GLN
1	C	230	ARG
1	A	167	SER
2	B	138	LEU
2	B	185	SER
1	C	58	ILE
1	C	130	ILE
2	B	115	HIS
1	C	226	LYS
2	D	124	GLN
2	D	165	LEU
1	A	184	GLN
2	B	181	GLU
1	C	138	PRO
1	C	59	HIS
1	C	109	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	175/175 (100%)	148 (85%)	27 (15%)	2	16
1	C	175/175 (100%)	154 (88%)	21 (12%)	5	23
2	B	68/72 (94%)	62 (91%)	6 (9%)	9	32
2	D	68/72 (94%)	56 (82%)	12 (18%)	2	11
All	All	486/494 (98%)	420 (86%)	66 (14%)	3	19

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LYS
1	A	68	THR
1	A	70	LEU

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Mol	Chain	Res	Type
1	A	73	ARG
1	A	75	ARG
1	A	85	ASP
1	A	88	THR
1	A	97	MET
1	A	103	VAL
1	A	108	ASP
1	A	127	LEU
1	A	130	ILE
1	A	136	GLU
1	A	137	ASN
1	A	141	ARG
1	A	146	ILE
1	A	151	MET
1	A	162	LEU
1	A	168	ASP
1	A	176	ILE
1	A	179	VAL
1	A	180	GLU
1	A	184	GLN
1	A	202	VAL
1	A	204	TRP
1	A	219	LEU
1	A	230	ARG
2	B	138	LEU
2	B	139	ASP
2	B	166	GLU
2	B	179	ASN
2	B	182	LYS
2	B	184	ILE
1	C	46	LYS
1	C	70	LEU
1	C	81	GLU
1	C	84	ILE
1	C	85	ASP
1	C	88	THR
1	C	93	ILE
1	C	95	SER
1	C	108	ASP
1	C	110	PHE
1	C	127	LEU
1	C	137	ASN

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Mol	Chain	Res	Type
1	C	141	ARG
1	C	146	ILE
1	C	155	VAL
1	C	162	LEU
1	C	176	ILE
1	C	187	GLN
1	C	214	CYS
1	C	219	LEU
1	C	230	ARG
2	D	119	LEU
2	D	125	SER
2	D	139	ASP
2	D	159	LEU
2	D	165	LEU
2	D	172	VAL
2	D	177	GLU
2	D	180	LEU
2	D	182	LYS
2	D	184	ILE
2	D	189	SER
2	D	191	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	59	HIS
1	A	86	GLN
1	A	137	ASN
1	A	140	ASN
1	A	187	GLN
1	A	212	ASN
2	B	118	GLN
2	B	129	GLN
2	B	179	ASN
1	C	56	GLN
1	C	59	HIS
1	C	137	ASN
1	C	157	ASN
1	C	182	ASN
1	C	184	GLN
1	C	212	ASN

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Mol	Chain	Res	Type
2	D	129	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	188/188 (100%)	-0.04	1 (0%) 87 68	67, 111, 204, 263	0
1	C	188/188 (100%)	-0.01	2 (1%) 78 54	71, 125, 209, 233	0
2	B	80/80 (100%)	-0.01	1 (1%) 75 50	54, 108, 224, 255	0
2	D	80/80 (100%)	-0.13	0 100 100	59, 113, 221, 266	0
All	All	536/536 (100%)	-0.04	4 (0%) 84 63	54, 116, 215, 266	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	92	PHE	2.7
1	C	61	ASN	2.4
2	B	192	PHE	2.3
1	C	92	PHE	2.1

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.