



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:38 PM UTC

PDB ID : 4GVR / pdb_00004gvr
Title : X-ray structure of the Archaeoglobus fulgidus methenyl-tetrahydromethanop
terin cyclohydrolase
Authors : Upadhyay, V.; Demmer, U.; Warkentin, E.; Moll, J.; Shima, S.; Ermler, U.
Deposited on : 2012-08-31
Resolution : 1.52 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

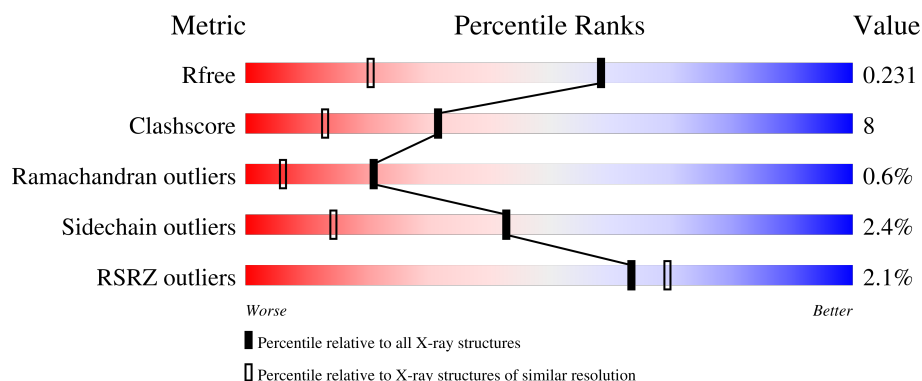
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.52 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	5890 (1.54-1.50)
Clashscore	190562	6116 (1.54-1.50)
Ramachandran outliers	187476	6002 (1.54-1.50)
Sidechain outliers	187428	5999 (1.54-1.50)
RSRZ outliers	180081	5891 (1.54-1.50)

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	316	 85% 13% .
1	B	316	 81% 15% . .
1	C	316	 83% 16% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8285 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 Methenyltetrahydromethanopterin cyclohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	10	0
			2513	1605	395	497	16			
1	B	315	Total	C	N	O	S	0	8	0
			2496	1594	391	493	18			
1	C	315	Total	C	N	O	S	0	12	0
			2536	1615	398	507	16			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	186	GLN	GLU	engineered mutation	UNP O28344
B	186	GLN	GLU	engineered mutation	UNP O28344
C	186	GLN	GLU	engineered mutation	UNP O28344

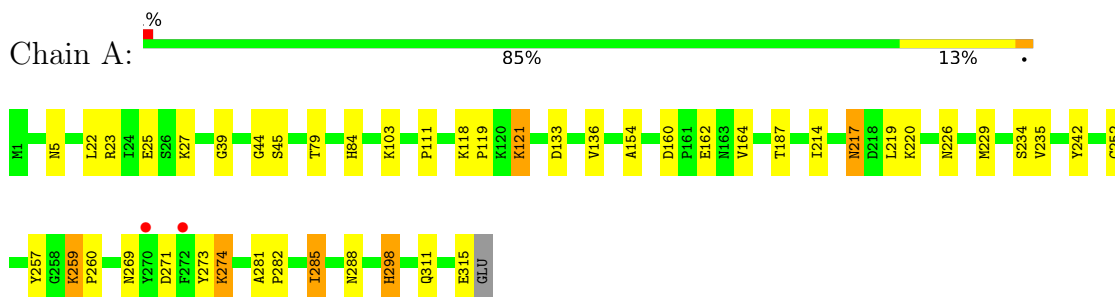
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	268	Total	O	0	0
			268	268		
2	B	208	Total	O	0	0
			208	208		
2	C	264	Total	O	0	0
			264	264		

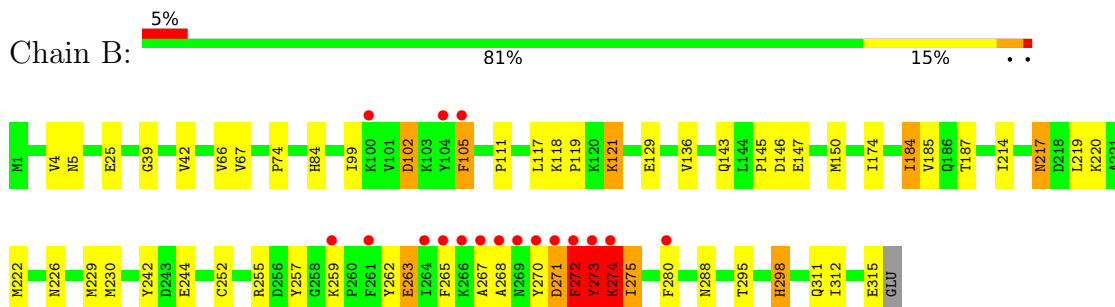
3 Residue-property plots [i](#)

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.

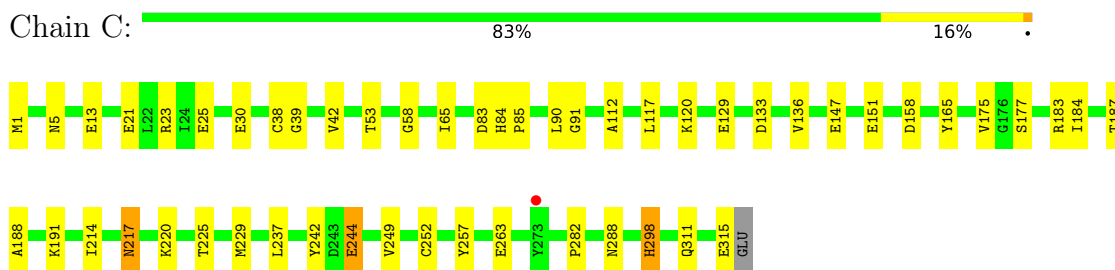
- Molecule 1: Methenyltetrahydromethanopterin cyclohydrolase



- Molecule 1: Methenyltetrahydromethanopterin cyclohydrolase



- Molecule 1: Methenyltetrahydromethanopterin cyclohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.54Å 122.08Å 79.39Å 90.00° 99.02° 90.00°	Depositor
Resolution (Å)	48.17 – 1.52 48.17 – 1.52	Depositor EDS
% Data completeness (in resolution range)	96.3 (48.17-1.52) 96.3 (48.17-1.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 1.52Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.164 , 0.231 0.163 , 0.231	Depositor DCC
R_{free} test set	8558 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.757	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8285	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.18	5/2560 (0.2%)	1.10	5/3476 (0.1%)
1	B	1.07	5/2543 (0.2%)	1.10	6/3452 (0.2%)
1	C	1.33	12/2583 (0.5%)	1.23	9/3507 (0.3%)
All	All	1.20	22/7686 (0.3%)	1.15	20/10435 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	112	ALA	N-CA	8.17	1.56	1.46
1	A	285	ILE	C-O	7.47	1.31	1.23
1	C	91	GLY	C-O	7.40	1.32	1.24
1	C	151	GLU	CD-OE1	7.02	1.38	1.25
1	A	22	LEU	C-O	6.54	1.32	1.24
1	B	263	GLU	CD-OE1	6.32	1.37	1.25
1	C	191	LYS	N-CA	6.23	1.53	1.46
1	C	298	HIS	ND1-CE1	6.12	1.38	1.32
1	C	58	GLY	N-CA	6.09	1.54	1.45
1	C	65	ILE	N-CA	6.05	1.53	1.46
1	B	263	GLU	CD-OE2	6.01	1.36	1.25
1	B	298	HIS	CE1-NE2	5.94	1.38	1.32
1	B	295	THR	C-O	5.93	1.30	1.24
1	A	27	LYS	C-O	5.83	1.30	1.23
1	C	225	THR	N-CA	5.76	1.53	1.46
1	A	45	SER	CA-C	5.68	1.59	1.52
1	C	237	LEU	N-CA	5.60	1.52	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	HIS	ND1-CE1	5.52	1.38	1.32
1	C	249	VAL	C-N	5.44	1.40	1.33
1	B	298	HIS	ND1-CE1	5.35	1.37	1.32
1	C	175	VAL	CA-CB	5.33	1.60	1.54
1	C	53	THR	CA-CB	5.10	1.61	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	274	LYS	N-CA-C	9.42	124.24	109.07
1	C	183	ARG	NE-CZ-NH2	-7.71	112.26	119.20
1	B	274	LYS	N-CA-C	7.09	125.89	110.80
1	B	271	ASP	N-CA-C	-6.24	97.50	110.80
1	C	90	LEU	O-C-N	5.78	130.77	122.28
1	C	183	ARG	NH1-CZ-NH2	5.54	126.50	119.30
1	C	249	VAL	CA-C-N	-5.50	114.49	119.94
1	C	249	VAL	C-N-CA	-5.50	114.49	119.94
1	C	188	ALA	N-CA-C	-5.38	105.33	111.14
1	B	275	ILE	N-CA-C	-5.26	100.87	108.71
1	C	38	CYS	N-CA-C	-5.25	106.10	112.88
1	C	177	SER	N-CA-C	5.25	116.69	111.07
1	A	84	HIS	CA-C-N	-5.23	113.53	119.28
1	A	84	HIS	C-N-CA	-5.23	113.53	119.28
1	B	273	TYR	N-CA-C	5.21	117.92	109.06
1	A	259	LYS	N-CA-C	5.16	117.16	110.40
1	B	185	VAL	N-CA-CB	5.12	116.54	110.55
1	C	85	PRO	N-CA-C	5.11	120.54	113.65
1	B	312	ILE	CB-CG1-CD1	5.08	124.46	113.80
1	A	103	LYS	N-CA-C	5.01	119.54	113.38

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	272	PHE	Peptide
1	B	273	TYR	Peptide
1	B	274	LYS	Peptide

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	2513	0	2479	37	0
1	B	2496	0	2461	67	0
1	C	2536	0	2483	34	0
2	A	268	0	0	6	0
2	B	208	0	0	5	0
2	C	264	0	0	5	0
All	All	8285	0	7423	122	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 8.

All (122) 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:273:TYR:HB2	1:B:275:ILE:N	1.56	1.20
1:B:273:TYR:CB	1:B:275:ILE:H	1.61	1.12
1:B:274:LYS:NZ	1:B:274:LYS:HB2	1.77	0.97
1:B:273:TYR:HB2	1:B:275:ILE:H	0.72	0.82
1:B:274:LYS:HB2	1:B:274:LYS:HZ2	1.45	0.81
1:B:25:GLU:HG2	1:B:42:VAL:HG11	1.63	0.80
1:B:145:PRO:CG	1:B:150[B]:MET:HE3	2.12	0.79
1:B:274:LYS:NZ	1:B:274:LYS:CB	2.49	0.76
1:A:288:ASN:HD21	1:B:214:ILE:H	1.34	0.76
1:B:217:ASN:ND2	1:B:220:LYS:H	1.84	0.76
1:A:214:ILE:H	1:C:288:ASN:HD21	1.35	0.75
1:B:121:LYS:HG3	2:B:492:HOH:O	1.87	0.74
1:A:160:ASP:OD1	1:A:162:GLU:HG3	1.89	0.73
1:A:315:GLU:OE2	1:C:298:HIS:HD2	1.71	0.72
1:B:145:PRO:HB2	1:B:150[B]:MET:HE3	1.72	0.72
1:B:270:TYR:CB	1:B:272:PHE:O	2.37	0.72
1:B:288:ASN:HD21	1:C:214:ILE:H	1.38	0.72
1:B:145:PRO:CB	1:B:150[B]:MET:HE3	2.21	0.71
1:B:268:ALA:O	1:B:270:TYR:N	2.23	0.70
1:B:265:PHE:CZ	1:B:271:ASP:HA	2.27	0.70
1:B:270:TYR:HB3	1:B:272:PHE:O	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:TYR:CG	1:B:272:PHE:O	2.45	0.69
1:B:117:LEU:HD21	1:B:129:GLU:HG3	1.77	0.66
1:C:187[A]:THR:HG21	1:C:229:MET:SD	2.36	0.65
1:B:187[A]:THR:HG21	1:B:229:MET:SD	2.37	0.65
1:B:242:TYR:OH	1:B:298:HIS:HE1	1.80	0.64
1:C:5:ASN:HD22	1:C:136:VAL:H	1.45	0.64
1:A:154:ALA:HB2	1:A:164:VAL:HG21	1.78	0.64
1:A:298:HIS:HD2	1:B:315:GLU:OE2	1.80	0.64
1:A:187[A]:THR:HG21	1:A:229:MET:SD	2.37	0.64
1:B:265:PHE:HE1	1:B:272:PHE:H	1.41	0.63
1:C:30:GLU:HG2	1:C:147:GLU:HG2	1.80	0.63
1:C:158:ASP:HB2	2:C:638:HOH:O	1.97	0.63
1:B:298:HIS:HD2	1:C:315:GLU:OE2	1.81	0.62
1:A:214:ILE:H	1:C:288:ASN:ND2	1.98	0.62
1:B:5:ASN:HD22	1:B:136:VAL:H	1.46	0.61
1:C:25:GLU:HG3	1:C:42:VAL:HG11	1.81	0.61
1:A:288:ASN:ND2	1:B:214:ILE:H	1.98	0.60
1:A:220:LYS:NZ	1:A:311:GLN:HE22	1.99	0.60
1:A:242:TYR:OH	1:A:298:HIS:HE1	1.85	0.59
1:A:23[B]:ARG:HD3	1:A:44:GLY:O	2.02	0.59
1:A:217:ASN:ND2	1:A:220:LYS:H	2.01	0.59
1:B:272:PHE:HE1	1:B:280:PHE:CZ	2.21	0.58
1:C:1:MET:HE2	2:C:547:HOH:O	2.03	0.58
1:C:217:ASN:ND2	1:C:220:LYS:H	2.02	0.58
1:B:25:GLU:HG2	1:B:42:VAL:CG1	2.35	0.56
1:B:230:MET:HG3	2:B:517:HOH:O	2.04	0.56
1:C:242:TYR:OH	1:C:298:HIS:HE1	1.89	0.56
1:A:23[B]:ARG:NH2	2:A:468:HOH:O	2.39	0.55
1:B:252:CYS:HA	1:B:257:TYR:CG	2.42	0.55
1:B:242:TYR:OH	1:B:298:HIS:CE1	2.60	0.55
1:B:4:VAL:HG23	1:B:111:PRO:HB3	1.88	0.54
1:C:217:ASN:HD21	1:C:220:LYS:H	1.55	0.54
1:C:220:LYS:NZ	1:C:311:GLN:HE22	2.04	0.54
1:C:220:LYS:HZ2	1:C:311:GLN:HE22	1.57	0.53
1:B:263:GLU:O	1:B:267:ALA:CB	2.55	0.53
1:B:230:MET:HE3	2:B:517:HOH:O	2.09	0.53
1:B:273:TYR:CD2	1:B:275:ILE:O	2.62	0.53
1:B:263:GLU:O	1:B:267:ALA:HB2	2.09	0.52
1:A:234:SER:HB3	1:B:74:PRO:HB3	1.92	0.52
1:B:274:LYS:CB	1:B:274:LYS:HZ3	2.22	0.52
2:A:662:HOH:O	1:C:84:HIS:HD2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ASN:HD22	1:A:217:ASN:C	2.20	0.49
1:B:5:ASN:ND2	1:B:136:VAL:H	2.08	0.49
1:C:187[B]:THR:HG23	1:C:282:PRO:HD3	1.95	0.49
1:B:145:PRO:HG2	1:B:150[B]:MET:HE3	1.93	0.49
1:A:242:TYR:OH	1:A:298:HIS:CE1	2.63	0.49
1:C:1:MET:CE	2:C:547:HOH:O	2.61	0.48
1:A:187[A]:THR:CG2	1:A:229:MET:SD	3.02	0.48
1:B:268:ALA:C	1:B:270:TYR:H	2.20	0.48
1:B:219:LEU:HD22	1:B:262:TYR:HE1	1.78	0.48
1:A:271:ASP:O	1:A:271:ASP:CG	2.54	0.48
1:B:298:HIS:CD2	1:C:315:GLU:OE2	2.66	0.48
1:A:217:ASN:HD21	1:A:220:LYS:H	1.61	0.48
1:A:23[B]:ARG:HD2	1:C:83:ASP:HB3	1.95	0.47
1:A:298:HIS:CD2	1:B:315:GLU:OE2	2.65	0.47
1:B:99:ILE:O	1:B:105:PHE:HA	2.14	0.47
1:A:5:ASN:HD22	1:A:136:VAL:H	1.63	0.47
1:B:226:ASN:HB3	1:B:230:MET:HE3	1.96	0.47
1:B:146:ASP:HB2	2:B:595:HOH:O	2.14	0.47
1:C:187[A]:THR:CG2	1:C:229:MET:SD	3.03	0.47
1:A:220:LYS:HZ2	1:A:311:GLN:HE22	1.63	0.46
1:A:281:ALA:HB1	1:A:282:PRO:HD2	1.97	0.46
1:B:220:LYS:NZ	1:B:311:GLN:HE22	2.14	0.46
1:C:184:ILE:O	1:C:187[B]:THR:HG22	2.15	0.46
1:B:66[B]:VAL:HG12	1:B:67:VAL:N	2.30	0.46
1:B:255:ARG:HH21	1:B:274:LYS:HG2	1.80	0.45
1:B:288:ASN:ND2	1:C:214:ILE:H	2.08	0.45
1:B:217:ASN:HD22	1:B:220:LYS:H	1.60	0.45
1:B:184:ILE:O	1:B:187[B]:THR:HG22	2.15	0.45
1:C:252:CYS:HA	1:C:257:TYR:CG	2.51	0.45
1:B:270:TYR:CD2	1:B:272:PHE:O	2.69	0.45
1:B:102:ASP:O	1:B:105:PHE:CE1	2.69	0.45
1:B:66[A]:VAL:HG13	1:B:66[A]:VAL:O	2.18	0.44
1:B:118:LYS:HA	1:B:119:PRO:HA	1.90	0.44
1:C:5:ASN:ND2	1:C:136:VAL:H	2.11	0.44
1:A:252:CYS:HA	1:A:257:TYR:CG	2.52	0.44
1:B:147:GLU:C	2:B:590:HOH:O	2.60	0.44
1:C:244:GLU:HG3	2:C:632:HOH:O	2.18	0.43
1:A:5:ASN:ND2	1:A:136:VAL:H	2.16	0.43
1:B:84:HIS:HE1	1:C:21:GLU:O	2.00	0.43
1:C:23[A]:ARG:NH1	2:C:661:HOH:O	2.51	0.43
1:A:226:ASN:ND2	2:A:577:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:LYS:HD2	1:B:311:GLN:HE22	1.83	0.43
1:A:79:THR:HG21	1:B:66[B]:VAL:HG13	1.99	0.43
1:A:315:GLU:OE2	1:C:298:HIS:CD2	2.60	0.43
1:B:265:PHE:HE1	1:B:272:PHE:N	2.14	0.43
1:B:217:ASN:HD21	1:B:220:LYS:H	1.65	0.42
1:A:269:ASN:HA	1:A:273:TYR:CE1	2.54	0.42
1:C:217:ASN:C	1:C:217:ASN:HD22	2.27	0.42
1:B:187[A]:THR:CG2	1:B:229:MET:SD	3.06	0.42
1:C:117:LEU:HD21	1:C:129:GLU:HG3	2.00	0.42
1:A:121:LYS:HD2	2:A:590:HOH:O	2.20	0.41
1:A:235:VAL:O	1:A:285:ILE:HA	2.20	0.41
1:C:13:GLU:OE2	1:C:165:TYR:OH	2.22	0.41
1:C:242:TYR:OH	1:C:298:HIS:CE1	2.70	0.40
1:A:25:GLU:HG2	2:A:476:HOH:O	2.20	0.40
1:A:259:LYS:O	1:A:260:PRO:C	2.63	0.40
2:A:565:HOH:O	1:B:174:ILE:HG12	2.21	0.40
1:A:5:ASN:ND2	1:A:111:PRO:HB2	2.36	0.40
1:A:118:LYS:HA	1:A:119:PRO:HA	1.96	0.40
1:B:272:PHE:HB3	1:B:273:TYR:HD1	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	323/316 (102%)	314 (97%)	8 (2%)	1 (0%)	36	18
1	B	321/316 (102%)	308 (96%)	9 (3%)	4 (1%)	10	1
1	C	325/316 (103%)	316 (97%)	8 (2%)	1 (0%)	36	18
All	All	969/948 (102%)	938 (97%)	25 (3%)	6 (1%)	21	6

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	102	ASP
1	A	39	GLY
1	B	274	LYS
1	B	39	GLY
1	C	39	GLY
1	B	272	PHE

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	275/266 (103%)	270 (98%)	5 (2%)	51	23
1	B	273/266 (103%)	263 (96%)	10 (4%)	30	6
1	C	277/266 (104%)	271 (98%)	6 (2%)	45	16
All	All	825/798 (103%)	804 (98%)	21 (2%)	43	13

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	LYS
1	A	133	ASP
1	A	217	ASN
1	A	219	LEU
1	A	274	LYS
1	B	105	PHE
1	B	121	LYS
1	B	143	GLN
1	B	184	ILE
1	B	217	ASN
1	B	222[A]	MET
1	B	222[B]	MET
1	B	244	GLU
1	B	259	LYS
1	B	274	LYS
1	C	120	LYS
1	C	133	ASP

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Mol	Chain	Res	Type
1	C	217	ASN
1	C	244	GLU
1	C	263[A]	GLU
1	C	263[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	98	GLN
1	A	143	GLN
1	A	217	ASN
1	A	226	ASN
1	A	288	ASN
1	A	298	HIS
1	A	311	GLN
1	B	5	ASN
1	B	98	GLN
1	B	217	ASN
1	B	288	ASN
1	B	298	HIS
1	B	311	GLN
1	C	5	ASN
1	C	84	HIS
1	C	98	GLN
1	C	217	ASN
1	C	288	ASN
1	C	298	HIS
1	C	311	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [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	315/316 (99%)	-0.43	2 (0%) 85 89	7, 20, 37, 58	10 (3%)
1	B	315/316 (99%)	0.03	17 (5%) 31 36	7, 25, 47, 79	8 (2%)
1	C	315/316 (99%)	-0.49	1 (0%) 90 92	7, 17, 37, 51	12 (3%)
All	All	945/948 (99%)	-0.30	20 (2%) 63 69	7, 20, 41, 79	30 (3%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	273	TYR	8.9
1	B	272	PHE	6.4
1	B	105	PHE	5.9
1	B	270	TYR	5.7
1	B	268	ALA	5.0
1	B	269	ASN	4.8
1	B	267	ALA	4.8
1	B	274	LYS	4.3
1	B	271	ASP	3.8
1	B	264	ILE	3.3
1	B	265	PHE	2.9
1	B	259	LYS	2.9
1	C	273	TYR	2.4
1	A	270	TYR	2.3
1	B	266	LYS	2.1
1	B	261	PHE	2.1
1	B	104	TYR	2.1
1	B	100	LYS	2.0
1	A	272	PHE	2.0
1	B	280	PHE	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.