



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:53 AM UTC

PDB ID : 3GYG / pdb_00003gyg
Title : Crystal structure of yhjK (haloacid dehalogenase-like hydrolase protein) from *Bacillus subtilis*
Authors : Nocek, B.; Stein, A.; Wu, R.; Jedrzejczak, R.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2009-04-03
Resolution : 2.45 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

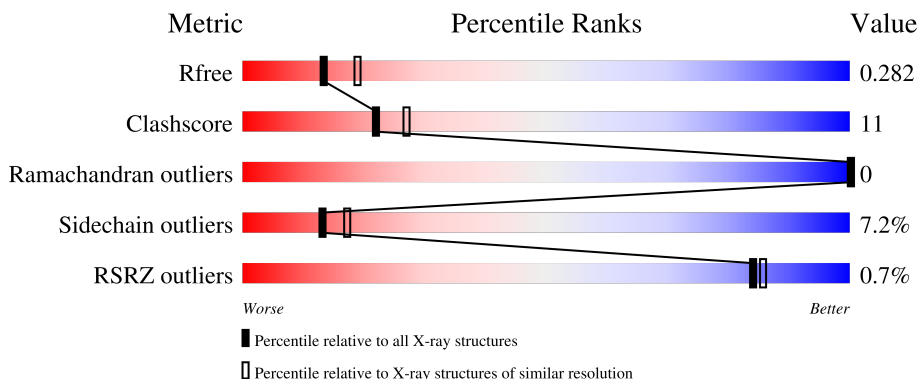
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	289	% 81% 17% ..
1	B	289	% 63% 16% • 20%
1	C	289	% 74% 19% • •
1	D	289	60% 23% • 13%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8445 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 NTD biosynthesis operon putative hydrolase ntdB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	Se	0	2	0
			2302	1458	383	453	3	5			
1	B	230	Total	C	N	O	S	Se	0	0	0
			1845	1172	304	362	2	5			
1	C	277	Total	C	N	O	S	Se	0	0	0
			2206	1398	363	439	3	3			
1	D	250	Total	C	N	O	S	Se	0	0	0
			2007	1279	330	390	3	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP O07565
A	-1	ASN	-	expression tag	UNP O07565
A	0	ALA	-	expression tag	UNP O07565
B	-2	SER	-	expression tag	UNP O07565
B	-1	ASN	-	expression tag	UNP O07565
B	0	ALA	-	expression tag	UNP O07565
C	-2	SER	-	expression tag	UNP O07565
C	-1	ASN	-	expression tag	UNP O07565
C	0	ALA	-	expression tag	UNP O07565
D	-2	SER	-	expression tag	UNP O07565
D	-1	ASN	-	expression tag	UNP O07565
D	0	ALA	-	expression tag	UNP O07565

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total 1	Mg 1	0	0
2	D	1	Total 1	Mg 1	0	0

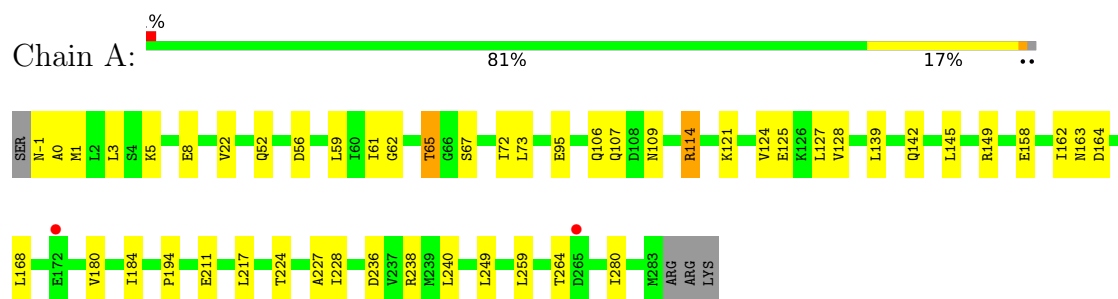
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	24	Total 24	O 24	0	0
3	B	15	Total 15	O 15	0	0
3	C	25	Total 25	O 25	0	0
3	D	17	Total 17	O 17	0	0

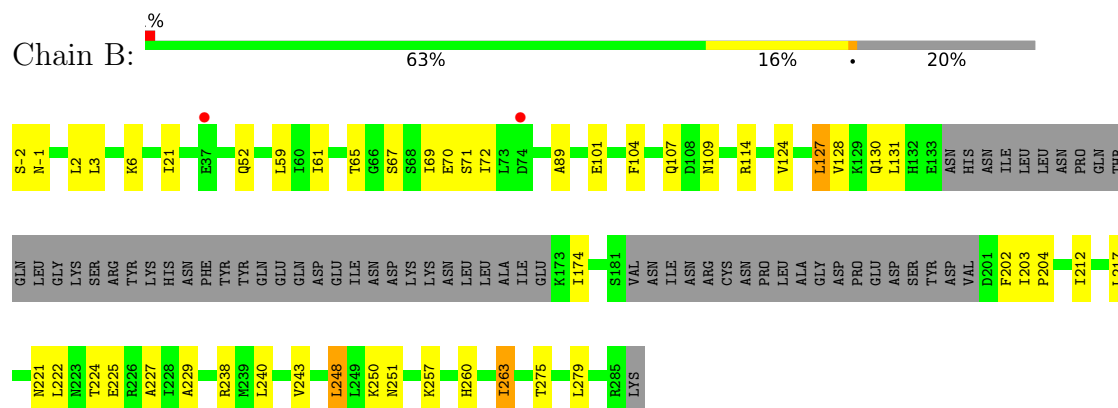
3 Residue-property plots

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.

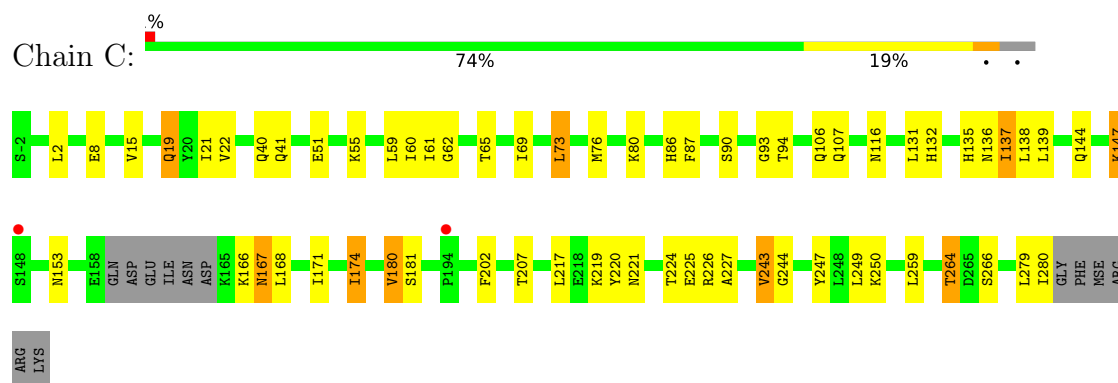
- Molecule 1: NTD biosynthesis operon putative hydrolase ntdB



- Molecule 1: NTD biosynthesis operon putative hydrolase ntdB

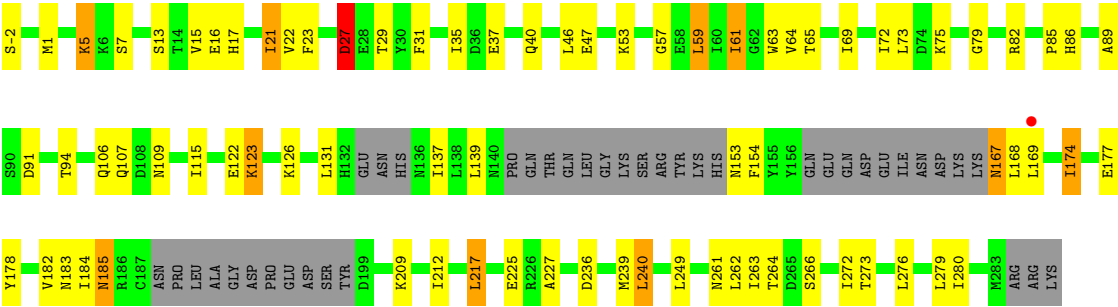


- Molecule 1: NTD biosynthesis operon putative hydrolase ntdB



- Molecule 1: NTD biosynthesis operon putative hydrolase ntdB

Chain D: 60% 23% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.08Å 111.73Å 135.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.45 40.00 – 2.45	Depositor EDS
% Data completeness (in resolution range)	93.4 (40.00-2.45) 93.4 (40.00-2.45)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0054	Depositor
R, R_{free}	0.210 , 0.277 0.224 , 0.282	Depositor DCC
R_{free} test set	1919 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8445	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 4.69% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.64	0/2350	0.79	0/3164
1	B	0.66	0/1875	0.80	0/2515
1	C	0.69	0/2247	0.82	0/3033
1	D	0.67	0/2037	0.89	2/2733 (0.1%)
All	All	0.66	0/8509	0.83	2/11445 (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	272	ILE	N-CA-C	-8.31	102.61	110.42
1	D	27	ASP	N-CA-C	5.09	118.24	110.20

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	2302	0	2210	37	0
1	B	1845	0	1775	43	0
1	C	2206	0	2086	53	0
1	D	2007	0	1945	55	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	24	0	0	0	0
3	B	15	0	0	0	0
3	C	25	0	0	0	0
3	D	17	0	0	1	0
All	All	8445	0	8016	175	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 11.

All (175) 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:114[A]:ARG:HG3	1:A:114[A]:ARG:NH1	1.56	1.09
1:D:65:THR:HG21	1:D:72:ILE:HD11	1.31	1.06
1:A:114[A]:ARG:HH11	1:A:114[A]:ARG:CG	1.69	1.05
1:D:22:VAL:HB	1:D:61:ILE:HD12	1.48	0.96
1:A:114[A]:ARG:HG3	1:A:114[A]:ARG:HH11	0.78	0.92
1:D:21:ILE:HD12	1:D:23:PHE:CZ	2.06	0.91
1:B:107:GLN:HE22	1:C:106:GLN:HA	1.38	0.88
1:D:21:ILE:HD12	1:D:23:PHE:CE2	2.10	0.86
1:B:65:THR:HG21	1:B:72:ILE:HD11	1.59	0.82
1:D:185:ASN:HD22	1:D:185:ASN:N	1.81	0.78
1:D:65:THR:CG2	1:D:72:ILE:HD11	2.14	0.76
1:B:131:LEU:CD1	1:B:174:ILE:HD13	2.18	0.74
1:A:1:MSE:HE3	1:A:73:LEU:HD12	1.71	0.71
1:C:137:ILE:HD12	1:C:167:ASN:ND2	2.05	0.70
1:B:65:THR:CG2	1:B:72:ILE:HD11	2.21	0.70
1:D:47:GLU:OE2	1:D:82:ARG:NH2	2.26	0.69
1:B:127:LEU:C	1:B:127:LEU:HD12	2.17	0.68
1:A:168:LEU:HD13	1:A:184:ILE:HG21	1.75	0.68
1:B:107:GLN:HE21	1:B:109:ASN:HD22	1.42	0.67
1:A:3:LEU:HD22	1:D:94:THR:HB	1.77	0.67
1:D:21:ILE:CD1	1:D:23:PHE:CZ	2.79	0.66
1:B:250:LYS:HA	1:B:263:ILE:HD11	1.76	0.66
1:C:249:LEU:HA	1:C:264:THR:HG22	1.79	0.65
1:C:227:ALA:CB	1:C:243:VAL:HG22	2.27	0.64
1:A:121:LYS:O	1:A:125:GLU:HG2	1.97	0.64
1:D:21:ILE:CD1	1:D:23:PHE:CE2	2.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:GLN:HE21	1:A:109:ASN:HD22	1.47	0.63
1:D:264:THR:HG22	1:D:266:SER:H	1.62	0.63
1:D:29:THR:HG23	1:D:249:LEU:HD11	1.81	0.62
1:A:128:VAL:HG22	1:A:139:LEU:HD22	1.82	0.60
1:D:53:LYS:HB3	1:D:59:LEU:HD23	1.85	0.59
1:D:22:VAL:CB	1:D:61:ILE:HD12	2.28	0.59
1:C:217:LEU:HD21	1:C:227:ALA:HB2	1.85	0.58
1:A:127:LEU:C	1:A:127:LEU:HD23	2.28	0.58
1:B:131:LEU:HD11	1:B:174:ILE:HG21	1.85	0.58
1:B:131:LEU:HD11	1:B:174:ILE:HD13	1.84	0.58
1:B:240:LEU:CD1	1:B:248:LEU:HD13	2.34	0.58
1:D:29:THR:HG23	1:D:249:LEU:CD1	2.35	0.57
1:A:1:MSE:HE3	1:A:73:LEU:CD1	2.35	0.57
1:D:16:GLU:HG2	1:D:17:HIS:CE1	2.39	0.57
1:C:167:ASN:HD22	1:C:167:ASN:C	2.13	0.57
1:C:19:GLN:O	1:C:226:ARG:HD2	2.06	0.56
1:C:59:LEU:HD12	1:C:59:LEU:C	2.31	0.56
1:B:59:LEU:C	1:B:59:LEU:HD12	2.31	0.55
1:B:127:LEU:HD12	1:B:128:VAL:N	2.21	0.55
1:C:2:LEU:C	1:C:2:LEU:HD23	2.32	0.55
1:D:139:LEU:HB3	1:D:154:PHE:HB3	1.88	0.55
1:D:13:SER:O	1:D:86:HIS:HE1	1.89	0.55
1:A:-1:ASN:HD21	1:A:1:MSE:HB3	1.70	0.55
1:C:131:LEU:HD23	1:C:174:ILE:HD11	1.89	0.55
1:B:6:LYS:NZ	1:C:116:ASN:HD22	2.04	0.54
1:B:21:ILE:HD12	1:B:217:LEU:HD23	1.88	0.54
1:B:21:ILE:HD11	1:B:222:LEU:HD12	1.90	0.54
1:D:131:LEU:HG	1:D:174:ILE:HD12	1.89	0.54
1:C:65:THR:O	1:C:90:SER:HA	2.07	0.53
1:C:264:THR:HG23	1:C:266:SER:O	2.08	0.53
1:A:52:GLN:NE2	1:A:56:ASP:OD1	2.41	0.53
1:C:227:ALA:HB1	1:C:243:VAL:CG2	2.39	0.53
1:C:224:THR:HG22	1:C:244:GLY:N	2.23	0.53
1:B:6:LYS:HZ2	1:C:116:ASN:HD22	1.55	0.53
1:B:275:THR:HG22	1:B:279:LEU:HD12	1.91	0.53
1:A:0:ALA:HB2	1:D:91:ASP:HB3	1.90	0.52
1:B:248:LEU:HD22	1:B:260:HIS:HB3	1.91	0.52
1:C:249:LEU:HA	1:C:264:THR:CG2	2.39	0.52
1:C:250:LYS:H	1:C:264:THR:HG22	1.73	0.52
1:A:107:GLN:HE22	1:D:106:GLN:HA	1.73	0.52
1:A:163:ASN:O	1:A:164:ASP:C	2.52	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:144:GLN:O	1:C:147:LYS:HB3	2.09	0.52
1:D:153:ASN:HB3	1:D:154:PHE:HD1	1.76	0.51
1:B:59:LEU:HD12	1:B:59:LEU:O	2.10	0.51
1:B:131:LEU:HD12	1:B:174:ILE:HD13	1.93	0.51
1:C:61:ILE:H	1:C:86:HIS:HD2	1.59	0.51
1:D:236:ASP:O	1:D:240:LEU:HD22	2.10	0.51
1:A:-1:ASN:ND2	1:A:1:MSE:HE2	2.26	0.50
1:D:27:ASP:OD2	1:D:27:ASP:N	2.41	0.50
1:A:5:LYS:NZ	1:A:8:GLU:OE1	2.42	0.50
1:A:95:GLU:HG3	1:D:-2:SER:O	2.11	0.50
1:A:106:GLN:HA	1:D:107:GLN:HE22	1.77	0.50
1:C:132:HIS:O	1:C:136:ASN:HA	2.12	0.50
1:C:264:THR:CG2	1:C:266:SER:O	2.60	0.50
1:D:123:LYS:NZ	1:D:178:TYR:O	2.37	0.50
1:D:183:ASN:CG	1:D:185:ASN:HD21	2.20	0.49
1:C:132:HIS:HD2	1:C:138:LEU:HD13	1.78	0.49
1:D:137:ILE:HD11	1:D:167:ASN:HB2	1.94	0.49
1:D:263:ILE:HG22	1:D:264:THR:O	2.13	0.48
1:D:35:ILE:HD12	1:D:79:GLY:HA2	1.95	0.48
1:B:21:ILE:CD1	1:B:222:LEU:HD12	2.44	0.47
1:B:257:LYS:HE2	1:B:263:ILE:HD13	1.95	0.47
1:D:31:PHE:CE2	1:D:75:LYS:HD3	2.49	0.47
1:D:185:ASN:N	1:D:185:ASN:ND2	2.51	0.47
1:C:153:ASN:ND2	1:C:202:PHE:H	2.13	0.47
1:C:40:GLN:NE2	1:C:80:LYS:HD2	2.28	0.47
1:A:61:ILE:HD12	1:A:62:GLY:N	2.30	0.47
1:B:250:LYS:HA	1:B:263:ILE:CD1	2.43	0.47
1:B:131:LEU:HD12	1:B:174:ILE:CD1	2.44	0.47
1:C:247:TYR:HE2	1:C:279:LEU:HD11	1.80	0.47
1:D:109:ASN:ND2	3:D:424:HOH:O	2.46	0.47
1:D:35:ILE:HG21	1:D:40:GLN:NE2	2.29	0.47
1:C:167:ASN:ND2	1:C:167:ASN:C	2.72	0.46
1:C:131:LEU:CD2	1:C:174:ILE:HD11	2.46	0.46
1:D:89:ALA:HB1	1:D:212:ILE:HD13	1.97	0.46
1:B:3:LEU:CD1	1:C:94:THR:HB	2.46	0.46
1:C:137:ILE:HD12	1:C:167:ASN:HD22	1.80	0.46
1:D:184:ILE:C	1:D:185:ASN:HD22	2.22	0.46
1:B:275:THR:HG22	1:B:279:LEU:CD1	2.45	0.46
1:D:122:GLU:HG2	1:D:126:LYS:HE2	1.98	0.46
1:A:-1:ASN:ND2	1:A:1:MSE:HB3	2.31	0.45
1:A:59:LEU:HD23	1:A:280:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:ALA:CB	1:B:243:VAL:HG21	2.45	0.45
1:B:227:ALA:HB1	1:B:243:VAL:HG12	1.98	0.45
1:A:22:VAL:HB	1:A:61:ILE:HD13	1.99	0.45
1:C:131:LEU:HG	1:C:174:ILE:HD12	1.99	0.45
1:C:227:ALA:HB3	1:C:243:VAL:HG22	1.97	0.45
1:B:127:LEU:C	1:B:127:LEU:CD1	2.87	0.45
1:C:61:ILE:HG22	1:C:86:HIS:HD2	1.82	0.45
1:C:73:LEU:O	1:C:76:MSE:HB2	2.17	0.45
1:D:29:THR:HA	1:D:249:LEU:HD12	1.99	0.45
1:B:65:THR:HG21	1:B:72:ILE:CD1	2.40	0.45
1:D:15:VAL:HG13	1:D:57:GLY:HA2	1.97	0.45
1:B:203:ILE:HG22	1:B:204:PRO:HD2	1.98	0.44
1:A:217:LEU:HD21	1:A:227:ALA:HB2	1.99	0.44
1:C:61:ILE:H	1:C:86:HIS:CD2	2.34	0.44
1:A:168:LEU:HD13	1:A:184:ILE:CG2	2.47	0.44
1:C:21:ILE:HD11	1:C:220:TYR:CD2	2.53	0.44
1:A:236:ASP:O	1:A:240:LEU:HG	2.17	0.44
1:B:131:LEU:CD1	1:B:174:ILE:CD1	2.92	0.44
1:C:21:ILE:HD13	1:C:60:ILE:HD11	1.99	0.44
1:A:65:THR:HG21	1:A:72:ILE:HD11	1.98	0.44
1:B:107:GLN:HE21	1:B:109:ASN:ND2	2.11	0.44
1:B:225:GLU:OE1	1:B:225:GLU:N	2.48	0.44
1:C:247:TYR:CE2	1:C:279:LEU:HD11	2.53	0.43
1:C:227:ALA:HB1	1:C:243:VAL:HG22	1.96	0.43
1:D:168:LEU:CD2	1:D:184:ILE:HD13	2.48	0.43
1:C:250:LYS:N	1:C:264:THR:HG22	2.33	0.43
1:A:124:VAL:HG11	1:D:7:SER:HB2	2.00	0.43
1:B:127:LEU:HD11	1:B:202:PHE:CE1	2.54	0.43
1:B:229:ALA:HB3	1:B:243:VAL:HG21	2.00	0.43
1:B:-2:SER:HB2	1:B:2:LEU:HD23	2.01	0.43
1:D:16:GLU:HG2	1:D:17:HIS:NE2	2.34	0.43
1:C:137:ILE:HD11	1:C:171:ILE:HD11	2.01	0.43
1:C:139:LEU:HD11	1:C:171:ILE:CD1	2.49	0.43
1:A:65:THR:CG2	1:A:72:ILE:HD11	2.49	0.42
1:B:114:ARG:HE	1:B:114:ARG:HB3	1.67	0.42
1:A:59:LEU:C	1:A:59:LEU:HD12	2.44	0.42
1:A:107:GLN:H	1:D:107:GLN:NE2	2.18	0.42
1:A:158:GLU:HB2	1:A:164:ASP:OD2	2.20	0.42
1:B:104:PHE:HB3	1:C:116:ASN:HD21	1.84	0.42
1:D:182:VAL:HG22	1:D:184:ILE:HG13	2.00	0.42
1:D:64:VAL:HG11	1:D:209:LYS:HG2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:LEU:HA	1:A:264:THR:OG1	2.19	0.42
1:D:63:TRP:CD2	1:D:85:PRO:HG3	2.55	0.42
1:B:89:ALA:HB1	1:B:212:ILE:HD13	2.02	0.42
1:C:59:LEU:HD12	1:C:59:LEU:O	2.20	0.42
1:D:46:LEU:HD13	1:D:273:THR:OG1	2.20	0.42
1:D:276:LEU:HD22	1:D:280:ILE:HD12	2.02	0.41
1:A:145:LEU:HD13	1:A:194:PRO:HG3	2.02	0.41
1:B:107:GLN:CB	1:C:107:GLN:HG2	2.50	0.41
1:C:93:GLY:HA3	1:C:207:THR:HB	2.02	0.41
1:D:167:ASN:OD1	1:D:167:ASN:N	2.53	0.41
1:C:147:LYS:O	1:C:147:LYS:HG3	2.14	0.41
1:D:5:LYS:HG2	1:D:73:LEU:HD22	2.03	0.41
1:D:53:LYS:CB	1:D:59:LEU:HD23	2.50	0.41
1:C:135:HIS:CD2	1:C:174:ILE:HD11	2.55	0.41
1:C:180:VAL:HG13	1:C:181:SER:N	2.36	0.41
1:D:217:LEU:HD11	1:D:227:ALA:HB2	2.02	0.41
1:C:51:GLU:OE2	1:C:55:LYS:HG3	2.21	0.41
1:A:22:VAL:HA	1:A:228:ILE:O	2.22	0.40
1:C:62:GLY:HA3	1:C:87:PHE:CZ	2.56	0.40
1:C:153:ASN:HD22	1:C:202:PHE:H	1.69	0.40
1:D:1:MSE:HE3	1:D:73:LEU:HD12	2.03	0.40
1:A:142:GLN:HE22	1:A:149:ARG:HE	1.69	0.40
1:B:124:VAL:O	1:B:128:VAL:HG23	2.21	0.40
1:B:251:ASN:H	1:B:251:ASN:ND2	2.20	0.40
1:C:139:LEU:HD11	1:C:171:ILE:HD11	2.04	0.40
1:D:264:THR:HG22	1:D:266:SER:N	2.32	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	285/289 (99%)	276 (97%)	9 (3%)	0	100	100
1	B	224/289 (78%)	216 (96%)	8 (4%)	0	100	100
1	C	273/289 (94%)	260 (95%)	13 (5%)	0	100	100
1	D	240/289 (83%)	229 (95%)	11 (5%)	0	100	100
All	All	1022/1156 (88%)	981 (96%)	41 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	253/257 (98%)	243 (96%)	10 (4%)	28	41
1	B	201/257 (78%)	186 (92%)	15 (8%)	12	16
1	C	240/257 (93%)	219 (91%)	21 (9%)	9	10
1	D	220/257 (86%)	199 (90%)	21 (10%)	8	8
All	All	914/1028 (89%)	847 (93%)	67 (7%)	13	17

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	THR
1	A	67	SER
1	A	114[A]	ARG
1	A	114[B]	ARG
1	A	162	ILE
1	A	180	VAL
1	A	211	GLU
1	A	224	THR
1	A	238	ARG
1	A	259	LEU
1	B	-1	ASN
1	B	52	GLN
1	B	61	ILE

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Mol	Chain	Res	Type
1	B	67	SER
1	B	69	ILE
1	B	70	GLU
1	B	71	SER
1	B	101	GLU
1	B	127	LEU
1	B	130	GLN
1	B	221	ASN
1	B	224	THR
1	B	238	ARG
1	B	248	LEU
1	B	263	ILE
1	C	8	GLU
1	C	15	VAL
1	C	19	GLN
1	C	22	VAL
1	C	41	GLN
1	C	69	ILE
1	C	73	LEU
1	C	137	ILE
1	C	147	LYS
1	C	166	LYS
1	C	167	ASN
1	C	168	LEU
1	C	174	ILE
1	C	180	VAL
1	C	219	LYS
1	C	221	ASN
1	C	225	GLU
1	C	243	VAL
1	C	259	LEU
1	C	264	THR
1	C	280	ILE
1	D	5	LYS
1	D	21	ILE
1	D	27	ASP
1	D	37	GLU
1	D	59	LEU
1	D	61	ILE
1	D	69	ILE
1	D	115	ILE
1	D	123	LYS

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Mol	Chain	Res	Type
1	D	167	ASN
1	D	169	LEU
1	D	174	ILE
1	D	177	GLU
1	D	185	ASN
1	D	217	LEU
1	D	225	GLU
1	D	239	MSE
1	D	240	LEU
1	D	261	ASN
1	D	262	LEU
1	D	279	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN
1	A	19	GLN
1	A	106	GLN
1	A	107	GLN
1	A	130	GLN
1	A	135	HIS
1	A	142	GLN
1	A	185	ASN
1	A	210	ASN
1	A	235	ASN
1	A	251	ASN
1	A	254	GLN
1	B	-1	ASN
1	B	107	GLN
1	B	109	ASN
1	B	210	ASN
1	B	221	ASN
1	B	251	ASN
1	B	254	GLN
1	C	40	GLN
1	C	52	GLN
1	C	86	HIS
1	C	116	ASN
1	C	132	HIS
1	C	135	HIS
1	C	144	GLN

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Mol	Chain	Res	Type
1	C	153	ASN
1	C	167	ASN
1	C	254	GLN
1	D	17	HIS
1	D	38	GLN
1	D	40	GLN
1	D	107	GLN
1	D	109	ASN
1	D	130	GLN
1	D	132	HIS
1	D	185	ASN
1	D	254	GLN
1	D	261	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	280/289 (96%)	-0.02	2 (0%) 84 86	7, 24, 35, 38	2 (0%)
1	B	225/289 (77%)	-0.13	2 (0%) 81 82	5, 22, 34, 43	0
1	C	273/289 (94%)	0.08	2 (0%) 84 86	15, 26, 37, 40	0
1	D	245/289 (84%)	0.00	1 (0%) 88 89	11, 23, 32, 36	0
All	All	1023/1156 (88%)	-0.01	7 (0%) 84 86	5, 24, 35, 43	2 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	172	GLU	2.8
1	A	265	ASP	2.6
1	C	148	SER	2.6
1	B	74	ASP	2.5
1	D	169	LEU	2.3
1	B	37	GLU	2.2
1	C	194	PRO	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	D	301	1/1	0.73	0.15	49,49,49,49	0
2	MG	C	301	1/1	0.81	0.15	49,49,49,49	0
2	MG	A	301	1/1	0.88	0.13	45,45,45,45	0
2	MG	B	301	1/1	0.91	0.09	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.