



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 12:30 AM EDT

PDB ID : 4I7U / pdb_00004i7u
Title : Dihydrodipicolinate Synthase of Agrobacterium tumefaciens
Authors : Atkinson, S.C.; Dogovski, C.; Dobson, R.C.J.; Perugini, M.A.
Deposited on : 2012-11-30
Resolution : 1.55 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

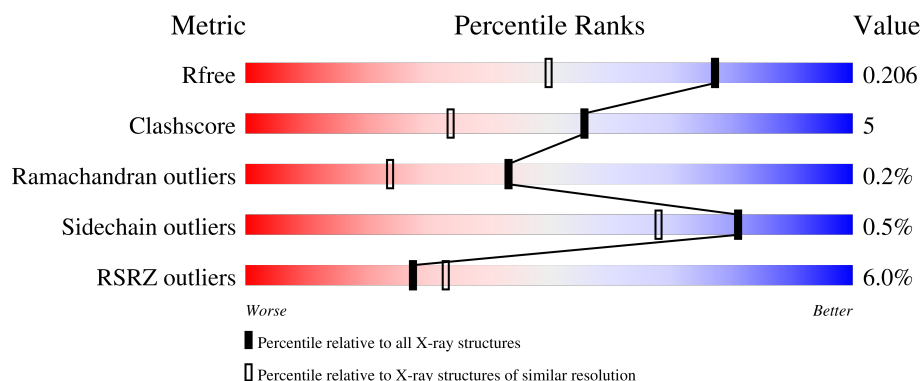
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	308	2% 88% 8% 5%
1	B	308	5% 89% 8% ..
1	C	308	5% 91% 5% .
1	D	308	12% 92% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10611 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 Dihydropicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	10	0
			2232	1407	381	430	14			
1	B	304	Total	C	N	O	S	0	19	0
			2355	1488	402	450	15			
1	C	295	Total	C	N	O	S	0	19	0
			2297	1452	398	432	15			
1	D	295	Total	C	N	O	S	0	10	0
			2240	1411	386	428	15			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP F0L8Z6
A	-12	ALA	-	expression tag	UNP F0L8Z6
A	-11	HIS	-	expression tag	UNP F0L8Z6
A	-10	HIS	-	expression tag	UNP F0L8Z6
A	-9	HIS	-	expression tag	UNP F0L8Z6
A	-8	HIS	-	expression tag	UNP F0L8Z6
A	-7	HIS	-	expression tag	UNP F0L8Z6
A	-6	HIS	-	expression tag	UNP F0L8Z6
A	-5	VAL	-	expression tag	UNP F0L8Z6
A	-4	ASP	-	expression tag	UNP F0L8Z6
A	-3	ASP	-	expression tag	UNP F0L8Z6
A	-2	ASP	-	expression tag	UNP F0L8Z6
A	-1	GLU	-	expression tag	UNP F0L8Z6
A	0	LYS	-	expression tag	UNP F0L8Z6
A	190	SER	GLY	SEE REMARK 999	UNP F0L8Z6
A	263	CYS	GLY	SEE REMARK 999	UNP F0L8Z6
B	-13	MET	-	expression tag	UNP F0L8Z6
B	-12	ALA	-	expression tag	UNP F0L8Z6
B	-11	HIS	-	expression tag	UNP F0L8Z6
B	-10	HIS	-	expression tag	UNP F0L8Z6
B	-9	HIS	-	expression tag	UNP F0L8Z6

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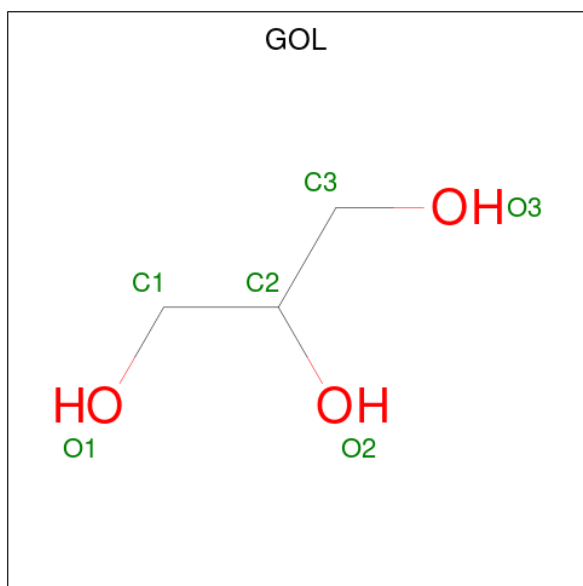
Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	HIS	-	expression tag	UNP F0L8Z6
B	-7	HIS	-	expression tag	UNP F0L8Z6
B	-6	HIS	-	expression tag	UNP F0L8Z6
B	-5	VAL	-	expression tag	UNP F0L8Z6
B	-4	ASP	-	expression tag	UNP F0L8Z6
B	-3	ASP	-	expression tag	UNP F0L8Z6
B	-2	ASP	-	expression tag	UNP F0L8Z6
B	-1	GLU	-	expression tag	UNP F0L8Z6
B	0	LYS	-	expression tag	UNP F0L8Z6
B	190	SER	GLY	SEE REMARK 999	UNP F0L8Z6
B	263	CYS	GLY	SEE REMARK 999	UNP F0L8Z6
C	-13	MET	-	expression tag	UNP F0L8Z6
C	-12	ALA	-	expression tag	UNP F0L8Z6
C	-11	HIS	-	expression tag	UNP F0L8Z6
C	-10	HIS	-	expression tag	UNP F0L8Z6
C	-9	HIS	-	expression tag	UNP F0L8Z6
C	-8	HIS	-	expression tag	UNP F0L8Z6
C	-7	HIS	-	expression tag	UNP F0L8Z6
C	-6	HIS	-	expression tag	UNP F0L8Z6
C	-5	VAL	-	expression tag	UNP F0L8Z6
C	-4	ASP	-	expression tag	UNP F0L8Z6
C	-3	ASP	-	expression tag	UNP F0L8Z6
C	-2	ASP	-	expression tag	UNP F0L8Z6
C	-1	GLU	-	expression tag	UNP F0L8Z6
C	0	LYS	-	expression tag	UNP F0L8Z6
C	190	SER	GLY	SEE REMARK 999	UNP F0L8Z6
C	263	CYS	GLY	SEE REMARK 999	UNP F0L8Z6
D	-13	MET	-	expression tag	UNP F0L8Z6
D	-12	ALA	-	expression tag	UNP F0L8Z6
D	-11	HIS	-	expression tag	UNP F0L8Z6
D	-10	HIS	-	expression tag	UNP F0L8Z6
D	-9	HIS	-	expression tag	UNP F0L8Z6
D	-8	HIS	-	expression tag	UNP F0L8Z6
D	-7	HIS	-	expression tag	UNP F0L8Z6
D	-6	HIS	-	expression tag	UNP F0L8Z6
D	-5	VAL	-	expression tag	UNP F0L8Z6
D	-4	ASP	-	expression tag	UNP F0L8Z6
D	-3	ASP	-	expression tag	UNP F0L8Z6
D	-2	ASP	-	expression tag	UNP F0L8Z6
D	-1	GLU	-	expression tag	UNP F0L8Z6
D	0	LYS	-	expression tag	UNP F0L8Z6
D	190	SER	GLY	SEE REMARK 999	UNP F0L8Z6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	263	CYS	GLY	SEE REMARK 999	UNP F0L8Z6

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	377	Total	O	0	0
			377	377		

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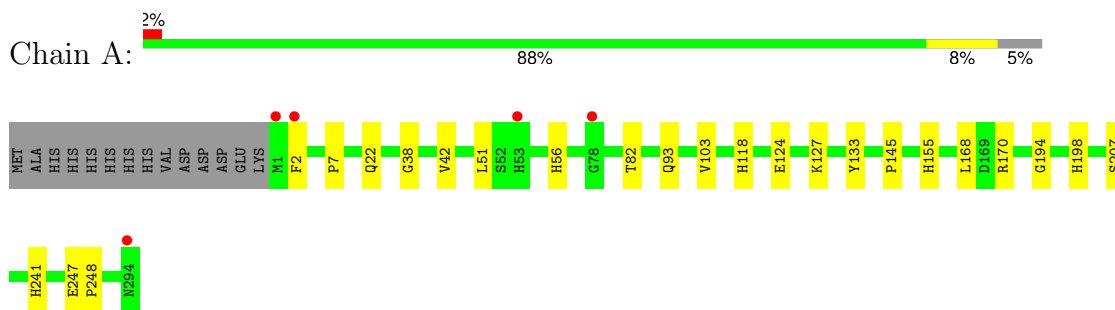
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	410	Total 410	O 410	0	0
3	C	363	Total 363	O 363	0	0
3	D	289	Total 289	O 289	0	0

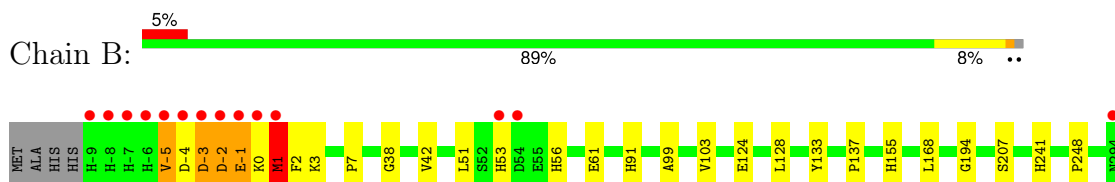
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

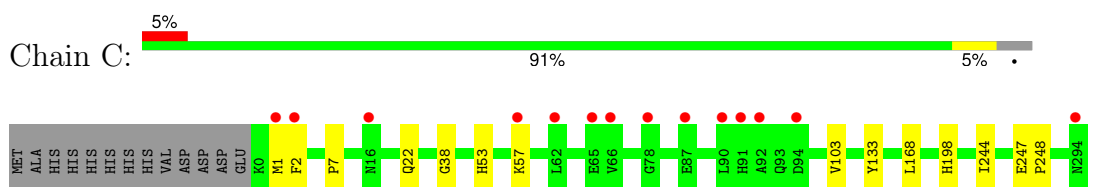
- Molecule 1: Dihydrodipicolinate synthase



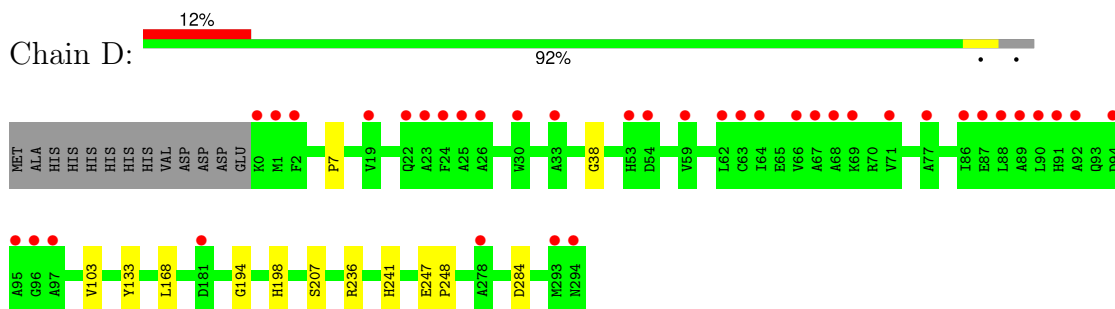
- Molecule 1: Dihydrodipicolinate synthase



- Molecule 1: Dihydrodipicolinate synthase



- Molecule 1: Dihydrodipicolinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.75Å 131.74Å 162.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	102.37 – 1.55 102.37 – 1.55	Depositor EDS
% Data completeness (in resolution range)	98.4 (102.37-1.55) 98.4 (102.37-1.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.189 , 0.204 0.195 , 0.206	Depositor DCC
R_{free} test set	14846 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10611	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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: GOL

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	0.46	1/2299 (0.0%)	0.74	0/3125
1	B	0.55	3/2459 (0.1%)	0.78	1/3341 (0.0%)
1	C	0.48	0/2397	0.74	0/3250
1	D	0.44	0/2312	0.75	0/3139
All	All	0.49	4/9467 (0.0%)	0.75	1/12855 (0.0%)

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	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	3	LYS	C-O	-5.62	1.17	1.23
1	A	2	PHE	C-O	-5.51	1.17	1.23
1	B	-3	ASP	N-CA	-5.11	1.40	1.46
1	B	2	PHE	C-O	-5.09	1.18	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	MET	N-CA-C	9.06	123.70	108.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	0	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	2232	0	2219	23	0
1	B	2355	0	2355	33	1
1	C	2297	0	2333	20	0
1	D	2240	0	2247	20	0
2	B	24	0	32	3	0
2	C	6	0	8	0	0
2	D	18	0	24	3	0
3	A	377	0	0	1	0
3	B	410	0	0	3	1
3	C	363	0	0	0	0
3	D	289	0	0	2	0
All	All	10611	0	9218	86	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (86) 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:168:LEU:HD12	1:D:168:LEU:HD12	1.16	1.13
1:D:236[B]:ARG:HH11	1:D:236[B]:ARG:HG3	1.04	1.08
1:B:168:LEU:HD12	1:C:168[A]:LEU:HD12	1.30	1.06
1:B:-3:ASP:O	1:B:-2:ASP:C	1.97	1.03
1:A:168:LEU:CD1	1:D:168:LEU:HD12	1.91	1.01
1:A:168:LEU:HD12	1:D:168:LEU:CD1	1.93	0.98
1:D:236[B]:ARG:HH11	1:D:236[B]:ARG:CG	1.77	0.98
1:B:248:PRO:HA	2:B:301:GOL:H2	1.52	0.91
1:B:168:LEU:CD1	1:C:168[A]:LEU:HD12	2.01	0.91
1:B:168:LEU:HD12	1:C:168[A]:LEU:CD1	2.00	0.90
1:D:236[B]:ARG:HG3	1:D:236[B]:ARG:NH1	1.87	0.88
1:A:93:GLN:HE22	1:A:127:LYS:H	1.17	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-2:ASP:O	1:B:-1:GLU:CB	2.30	0.80
1:B:-4:ASP:OD1	1:B:-4:ASP:C	2.34	0.70
1:C:103:VAL:HA	1:C:133[A]:TYR:HB3	1.76	0.66
1:C:1:MET:HG2	1:C:2:PHE:CD1	2.32	0.64
1:D:103:VAL:HA	1:D:133[B]:TYR:HB3	1.79	0.64
1:A:198:HIS:HE1	1:D:194:GLY:H	1.46	0.62
1:A:93:GLN:NE2	1:A:127:LYS:H	1.94	0.61
1:A:194:GLY:H	1:D:198:HIS:HE1	1.48	0.61
1:B:194:GLY:H	1:C:198:HIS:HE1	1.48	0.61
1:B:42[B]:VAL:HG13	1:B:51:LEU:HD12	1.85	0.59
1:A:194:GLY:H	1:D:198:HIS:CE1	2.20	0.59
1:A:42[A]:VAL:HG23	1:A:51:LEU:HD12	1.85	0.59
1:A:103:VAL:HA	1:A:133[B]:TYR:HB3	1.85	0.59
1:B:-3:ASP:O	1:B:-1:GLU:N	2.35	0.58
1:B:194:GLY:H	1:C:198:HIS:CE1	2.21	0.58
1:A:198:HIS:CE1	1:D:194:GLY:H	2.22	0.58
1:A:82:THR:OG1	1:A:118:HIS:HD2	1.87	0.57
1:D:236[B]:ARG:CG	1:D:236[B]:ARG:NH1	2.47	0.57
1:B:-5:VAL:HG21	1:B:1:MET:HB2	1.86	0.57
1:A:207:SER:OG	1:A:241:HIS:HD2	1.87	0.57
1:B:103:VAL:HA	1:B:133[B]:TYR:HB3	1.85	0.57
2:B:301:GOL:H11	3:B:542:HOH:O	2.04	0.57
1:A:7:PRO:HD2	1:A:38:GLY:O	2.04	0.57
1:B:61:GLU:OE2	1:B:91:HIS:HE1	1.87	0.57
1:D:207:SER:OG	1:D:241:HIS:HD2	1.88	0.57
1:D:133[A]:TYR:OH	2:D:301:GOL:H31	2.06	0.56
1:B:137:PRO:HA	2:B:302:GOL:H2	1.88	0.56
1:A:103:VAL:HA	1:A:133[A]:TYR:HB3	1.88	0.56
1:B:207:SER:OG	1:B:241:HIS:HD2	1.88	0.56
1:D:103:VAL:HA	1:D:133[A]:TYR:HB3	1.89	0.55
1:B:241:HIS:HE1	3:B:444:HOH:O	1.91	0.53
1:B:-3:ASP:O	1:B:-2:ASP:O	2.26	0.52
1:C:103:VAL:HA	1:C:133[B]:TYR:HB3	1.92	0.52
1:C:1:MET:SD	1:C:2:PHE:CE1	3.04	0.51
1:B:-5:VAL:O	1:B:-2:ASP:N	2.44	0.51
1:A:241:HIS:HE1	3:A:334:HOH:O	1.93	0.51
1:B:103:VAL:HA	1:B:133[A]:TYR:HB3	1.92	0.51
2:D:303:GOL:H32	3:D:505:HOH:O	2.11	0.50
1:B:-3:ASP:O	1:B:-1:GLU:CB	2.59	0.49
1:D:7:PRO:HD2	1:D:38:GLY:O	2.12	0.49
1:B:-4:ASP:OD1	1:B:-4:ASP:O	2.30	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42[A]:VAL:HG22	1:A:42[A]:VAL:O	2.13	0.48
1:D:241:HIS:HE1	3:D:441:HOH:O	1.95	0.48
1:B:124[B]:GLU:HG2	1:B:155:HIS:HE1	1.78	0.47
1:C:7:PRO:HD2	1:C:38:GLY:O	2.14	0.47
1:B:-3:ASP:O	1:B:-3:ASP:OD2	2.32	0.47
1:B:-5:VAL:C	1:B:-3:ASP:H	2.23	0.46
1:D:247:GLU:HB3	1:D:248:PRO:HD2	1.96	0.46
1:B:7:PRO:HD2	1:B:38:GLY:O	2.16	0.46
1:C:1:MET:HE3	1:C:1:MET:HB2	1.68	0.46
1:C:247:GLU:HB3	1:C:248:PRO:CD	2.45	0.46
1:D:247:GLU:HB3	1:D:248:PRO:CD	2.46	0.46
1:D:284:ASP:OD2	2:D:303:GOL:H12	2.17	0.45
1:B:99:ALA:HA	1:B:128[B]:LEU:HG	1.98	0.45
1:A:42[A]:VAL:HG21	1:A:56:HIS:CD2	2.51	0.45
1:C:22[A]:GLN:H	1:C:22[A]:GLN:CD	2.26	0.44
1:B:91:HIS:HD2	3:B:503:HOH:O	2.01	0.44
1:C:2:PHE:CD1	1:C:2:PHE:N	2.86	0.44
1:A:42[A]:VAL:O	1:A:42[A]:VAL:HG13	2.17	0.44
1:A:247:GLU:HB3	1:A:248:PRO:CD	2.48	0.44
1:B:42[B]:VAL:HG11	1:B:56:HIS:CD2	2.52	0.44
1:C:53:HIS:O	1:C:57:LYS:HG3	2.17	0.44
1:A:145:PRO:HG3	1:A:170:ARG:HD2	2.01	0.43
1:C:1:MET:SD	1:C:2:PHE:CZ	3.11	0.43
1:C:1:MET:HG2	1:C:2:PHE:CG	2.53	0.42
1:B:-5:VAL:HG11	1:B:1:MET:HB2	2.02	0.42
1:C:247:GLU:HB3	1:C:248:PRO:HD2	2.01	0.42
1:B:124[B]:GLU:HG2	1:B:155:HIS:CE1	2.54	0.41
1:C:1:MET:HG2	1:C:2:PHE:CE1	2.56	0.41
1:C:168[B]:LEU:HD13	1:C:168[B]:LEU:HA	1.87	0.41
1:B:124[A]:GLU:HG3	1:B:155:HIS:HE1	1.86	0.41
1:A:124:GLU:HG3	1:A:155:HIS:HE1	1.85	0.40
1:A:124:GLU:HG3	1:A:155:HIS:CE1	2.56	0.40
1:B:-5:VAL:C	1:B:-3:ASP:N	2.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-2:ASP:O	3:B:698:HOH:O[4_545]	1.71	0.49

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	302/308 (98%)	296 (98%)	6 (2%)	0	100	100
1	B	321/308 (104%)	313 (98%)	6 (2%)	2 (1%)	21	6
1	C	312/308 (101%)	306 (98%)	6 (2%)	0	100	100
1	D	303/308 (98%)	298 (98%)	5 (2%)	0	100	100
All	All	1238/1232 (100%)	1213 (98%)	23 (2%)	2 (0%)	43	24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	-1	GLU
1	B	-2	ASP

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	236/243 (97%)	234 (99%)	2 (1%)	73	56
1	B	253/243 (104%)	250 (99%)	3 (1%)	63	40
1	C	246/243 (101%)	244 (99%)	2 (1%)	73	56
1	D	238/243 (98%)	238 (100%)	0	100	100
All	All	973/972 (100%)	966 (99%)	7 (1%)	81	59

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22[A]	GLN
1	A	22[B]	GLN
1	B	-5	VAL
1	B	1	MET
1	B	53	HIS
1	C	244[A]	ILE
1	C	244[B]	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	GLN
1	A	108	ASN
1	A	118	HIS
1	A	196	ASN
1	A	198	HIS
1	A	241	HIS
1	B	56	HIS
1	B	91	HIS
1	B	108	ASN
1	B	196	ASN
1	B	241	HIS
1	C	108	ASN
1	C	196	ASN
1	C	198	HIS
1	D	196	ASN
1	D	198	HIS
1	D	241	HIS

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

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	D	303	-	5,5,5	0.41	0	5,5,5	0.21	0
2	GOL	D	301	-	5,5,5	0.32	0	5,5,5	0.38	0
2	GOL	B	303	-	5,5,5	0.38	0	5,5,5	0.32	0
2	GOL	B	302	-	5,5,5	0.35	0	5,5,5	0.29	0
2	GOL	C	301	-	5,5,5	0.40	0	5,5,5	0.34	0
2	GOL	B	301	-	5,5,5	0.45	0	5,5,5	0.68	0
2	GOL	D	302	-	5,5,5	0.35	0	5,5,5	0.64	0
2	GOL	B	304	-	5,5,5	0.38	0	5,5,5	0.37	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	D	303	-	-	4/4/4/4	-
2	GOL	D	301	-	-	0/4/4/4	-
2	GOL	B	303	-	-	1/4/4/4	-
2	GOL	B	302	-	-	2/4/4/4	-
2	GOL	C	301	-	-	3/4/4/4	-
2	GOL	B	301	-	-	2/4/4/4	-
2	GOL	D	302	-	-	0/4/4/4	-
2	GOL	B	304	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	302	GOL	C1-C2-C3-O3
2	B	302	GOL	O2-C2-C3-O3
2	B	304	GOL	C1-C2-C3-O3
2	B	304	GOL	O2-C2-C3-O3
2	C	301	GOL	C1-C2-C3-O3
2	D	303	GOL	O1-C1-C2-O2
2	D	303	GOL	O1-C1-C2-C3
2	D	303	GOL	C1-C2-C3-O3
2	B	301	GOL	C1-C2-C3-O3
2	C	301	GOL	O1-C1-C2-C3
2	B	301	GOL	O2-C2-C3-O3
2	D	303	GOL	O2-C2-C3-O3
2	B	304	GOL	O1-C1-C2-O2
2	C	301	GOL	O2-C2-C3-O3
2	B	303	GOL	O2-C2-C3-O3

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	303	GOL	2	0
2	D	301	GOL	1	0
2	B	302	GOL	1	0
2	B	301	GOL	2	0

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	294/308 (95%)	0.05	5 (1%) 69 77	6, 17, 24, 32	10 (3%)
1	B	304/308 (98%)	0.10	14 (4%) 37 45	6, 14, 24, 52	19 (6%)
1	C	295/308 (95%)	0.38	14 (4%) 36 44	9, 18, 28, 33	19 (6%)
1	D	295/308 (95%)	0.91	38 (12%) 7 9	11, 22, 33, 37	10 (3%)
All	All	1188/1232 (96%)	0.36	71 (5%) 27 33	6, 18, 30, 52	58 (4%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	-2	ASP	13.7
1	B	-3	ASP	11.4
1	B	-1	GLU	8.5
1	B	0	LYS	8.3
1	B	-9	HIS	8.0
1	B	-5	VAL	7.8
1	B	-7[A]	HIS	6.3
1	B	-4	ASP	5.8
1	C	2	PHE	5.7
1	C	1	MET	5.6
1	D	2	PHE	5.0
1	D	1	MET	4.8
1	C	91[A]	HIS	4.5
1	D	91	HIS	3.9
1	B	-6	HIS	3.7
1	B	-8	HIS	3.6
1	B	1	MET	3.5
1	D	92	ALA	3.3
1	D	22	GLN	3.2
1	C	57	LYS	3.2
1	D	19	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	90	LEU	3.1
1	D	87	GLU	3.0
1	D	89	ALA	3.0
1	D	97	ALA	3.0
1	D	26	ALA	2.9
1	D	294	ASN	2.9
1	C	78	GLY	2.9
1	A	294	ASN	2.8
1	C	65	GLU	2.8
1	D	25	ALA	2.8
1	B	53	HIS	2.8
1	A	1	MET	2.8
1	D	66	VAL	2.7
1	C	94[A]	ASP	2.7
1	D	62	LEU	2.7
1	C	66	VAL	2.6
1	C	294	ASN	2.6
1	D	64	ILE	2.6
1	D	68	ALA	2.6
1	A	2	PHE	2.6
1	D	69	LYS	2.6
1	D	30	TRP	2.5
1	C	90	LEU	2.5
1	D	53	HIS	2.5
1	C	87	GLU	2.5
1	D	59	VAL	2.5
1	C	92	ALA	2.5
1	D	94[A]	ASP	2.5
1	D	54	ASP	2.5
1	D	96	GLY	2.4
1	D	67	ALA	2.4
1	B	294	ASN	2.4
1	D	24	PHE	2.3
1	D	0	LYS	2.3
1	D	23	ALA	2.3
1	D	293	MET	2.3
1	C	16	ASN	2.3
1	D	71	VAL	2.3
1	C	62	LEU	2.2
1	D	77	ALA	2.2
1	D	88	LEU	2.1
1	D	33	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	95	ALA	2.1
1	D	278	ALA	2.1
1	B	54	ASP	2.1
1	D	181[A]	ASP	2.1
1	A	78	GLY	2.1
1	D	86	ILE	2.1
1	A	53	HIS	2.0
1	D	63	CYS	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(Å ²)	Q<0.9
2	GOL	B	303	6/6	0.71	0.27	58,58,58,59	0
2	GOL	B	304	6/6	0.71	0.27	56,56,56,56	0
2	GOL	C	301	6/6	0.73	0.25	37,39,39,39	0
2	GOL	D	301	6/6	0.75	0.22	33,34,35,35	0
2	GOL	B	301	6/6	0.76	0.20	22,28,28,31	0
2	GOL	D	302	6/6	0.77	0.20	32,34,35,36	0
2	GOL	D	303	6/6	0.77	0.19	38,39,40,42	0
2	GOL	B	302	6/6	0.78	0.19	38,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.