



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:10 AM UTC

PDB ID : 2GAC / pdb_00002gac
Title : T152C MUTANT GLYCOSYLASPARAGINASE FROM FLAVOBACTERIUM MENINGOSEPTICUM
Authors : Guo, H.-C.; Xu, Q.
Deposited on : 1998-05-29
Resolution : 2.10 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

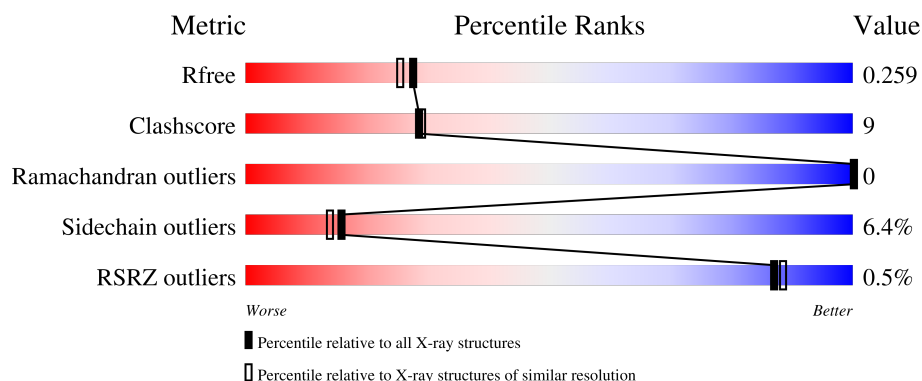
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	151	
1	C	151	
2	B	144	
2	D	144	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4432 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 GLYCOSYLASPARAGINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	136	Total	C	N	O	S	0	0	0
			1056	665	184	201	6			
1	C	136	Total	C	N	O	S	0	0	0
			1056	665	184	201	6			

- Molecule 2 is a protein called GLYCOSYLASPARAGINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	139	Total	C	N	O	S	0	0	0
			1037	642	191	196	8			
2	D	139	Total	C	N	O	S	0	0	0
			1037	642	191	196	8			

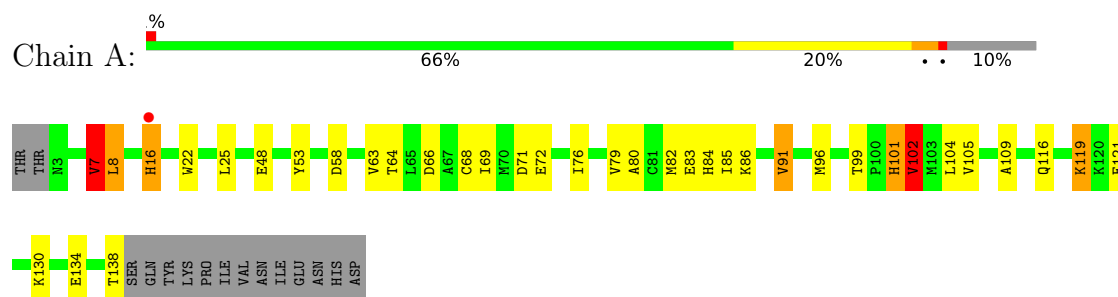
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		
3	B	52	Total	O	0	0
			52	52		
3	C	71	Total	O	0	0
			71	71		
3	D	54	Total	O	0	0
			54	54		

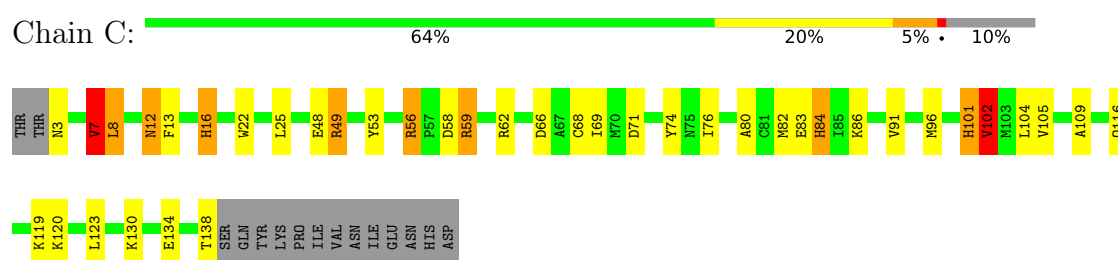
3 Residue-property plots

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.

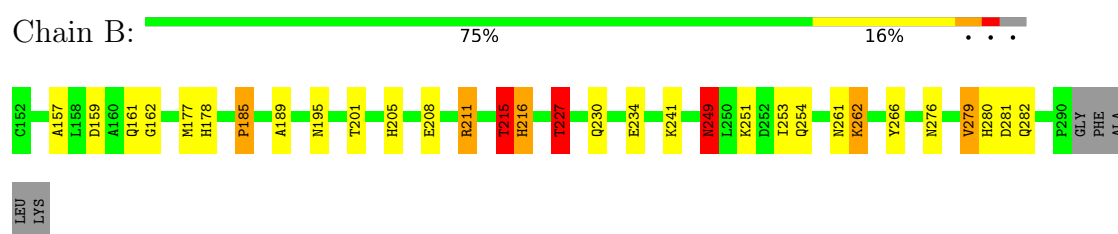
• Molecule 1: GLYCOSYLASPARAGINASE



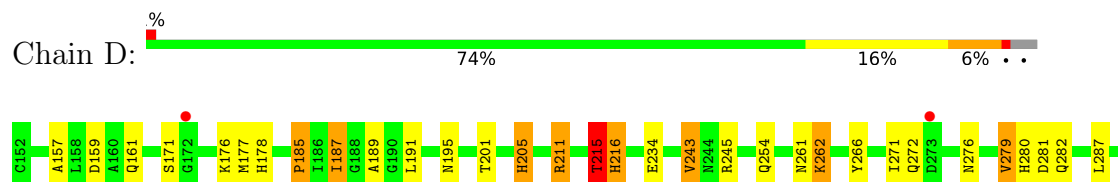
• Molecule 1: GLYCOSYLASPARAGINASE



• Molecule 2: GLYCOSYLASPARAGINASE



• Molecule 2: GLYCOSYLASPARAGINASE



P250
GLY
PHE
ALA
LEU
LYS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.20Å 97.30Å 61.80Å 90.00° 90.30° 90.00°	Depositor
Resolution (Å)	6.00 – 2.10 6.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.5 (6.00-2.10) 91.7 (6.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.10 (at 2.10Å)	Xtriage
Refinement program	XTALVIEW, X-PLOR 3.1	Depositor
R, R_{free}	0.233 , 0.280 0.220 , 0.259	Depositor DCC
R_{free} test set	3058 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.624	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.87 , 115.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4432	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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	0.89	4/1076 (0.4%)	1.68	16/1453 (1.1%)
1	C	0.91	3/1076 (0.3%)	1.77	23/1453 (1.6%)
2	B	0.84	4/1050 (0.4%)	1.63	19/1412 (1.3%)
2	D	0.88	6/1050 (0.6%)	1.69	21/1412 (1.5%)
All	All	0.88	17/4252 (0.4%)	1.69	79/5730 (1.4%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	178	HIS	CD2-NE2	-7.24	1.29	1.37
1	C	84	HIS	CD2-NE2	-7.04	1.30	1.37
1	C	101	HIS	CD2-NE2	-6.79	1.30	1.37
2	B	178	HIS	CD2-NE2	-6.77	1.30	1.37
1	A	101	HIS	CD2-NE2	-6.76	1.30	1.37
2	B	205	HIS	CD2-NE2	-6.73	1.30	1.37
2	D	216	HIS	CD2-NE2	-6.64	1.30	1.37
1	A	84	HIS	CD2-NE2	-6.55	1.30	1.37
1	C	16	HIS	CD2-NE2	-6.50	1.30	1.37
2	B	216	HIS	CD2-NE2	-6.46	1.30	1.37
2	B	280	HIS	CD2-NE2	-6.14	1.31	1.37
2	D	280	HIS	CD2-NE2	-6.09	1.31	1.37
1	A	16	HIS	CD2-NE2	-6.04	1.31	1.37
2	D	243	VAL	CA-CB	5.78	1.62	1.54
2	D	205	HIS	CD2-NE2	-5.74	1.31	1.37
1	A	84	HIS	CG-ND1	-5.67	1.32	1.38
2	D	205	HIS	CG-ND1	-5.07	1.32	1.38

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	215	THR	N-CA-CB	-10.72	93.66	110.28
2	D	215	THR	N-CA-CB	-10.63	93.81	110.28
1	C	102	VAL	N-CA-CB	-9.95	98.69	112.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	187	ILE	O-C-N	-9.59	112.56	121.87
1	C	102	VAL	CB-CA-C	9.52	119.98	110.65
1	A	102	VAL	CB-CA-C	9.35	119.42	110.63
1	A	58	ASP	CA-CB-CG	9.32	121.92	112.60
1	C	59	ARG	CB-CG-CD	-8.92	90.79	111.30
2	D	171	SER	N-CA-C	-8.50	96.81	110.32
2	B	211	ARG	NE-CZ-NH2	-8.48	111.57	119.20
2	D	211	ARG	NE-CZ-NH2	-8.35	111.69	119.20
1	A	16	HIS	ND1-CG-CD2	8.27	114.37	106.10
1	A	71	ASP	CA-CB-CG	8.21	120.81	112.60
1	C	7	VAL	CB-CA-C	-8.16	98.47	110.62
2	D	187	ILE	CA-C-O	-8.04	112.59	120.95
1	A	7	VAL	CB-CA-C	-8.04	98.65	110.62
1	A	102	VAL	N-CA-CB	-7.99	99.26	111.57
1	C	71	ASP	CA-CB-CG	7.95	120.55	112.60
1	C	58	ASP	CA-CB-CG	7.69	120.29	112.60
1	C	16	HIS	ND1-CG-CD2	7.55	113.65	106.10
1	C	48	GLU	CA-C-O	7.54	128.76	120.92
2	B	280	HIS	CB-CG-CD2	-7.49	121.46	131.20
2	D	282	GLN	OE1-CD-NE2	-7.30	115.30	122.60
2	D	171	SER	O-C-N	-7.26	114.93	123.06
2	D	211	ARG	CA-CB-CG	7.19	128.47	114.10
2	B	211	ARG	CA-CB-CG	7.15	128.41	114.10
2	B	249	ASN	OD1-CG-ND2	-7.10	115.50	122.60
2	D	280	HIS	CB-CG-CD2	-7.09	121.98	131.20
1	C	59	ARG	NE-CZ-NH2	-6.88	113.00	119.20
2	D	261	ASN	CA-CB-CG	6.82	119.42	112.60
2	B	159	ASP	CA-CB-CG	6.76	119.36	112.60
2	D	276	ASN	OD1-CG-ND2	-6.68	115.92	122.60
2	D	215	THR	CB-CA-C	6.64	123.42	110.67
2	B	215	THR	CB-CA-C	6.63	123.41	110.67
2	B	276	ASN	OD1-CG-ND2	-6.62	115.98	122.60
1	C	13	PHE	CA-CB-CG	6.62	120.42	113.80
1	C	12	ASN	OD1-CG-ND2	-6.59	116.01	122.60
1	C	66	ASP	CA-CB-CG	6.45	119.05	112.60
2	D	281	ASP	CA-CB-CG	6.34	118.94	112.60
2	B	178	HIS	CB-CG-CD2	-6.31	123.00	131.20
2	B	261	ASN	CA-CB-CG	6.30	118.90	112.60
2	D	245	ARG	CA-CB-CG	6.24	126.57	114.10
1	A	66	ASP	CA-CB-CG	6.21	118.81	112.60
2	D	159	ASP	CA-CB-CG	6.14	118.75	112.60
1	A	16	HIS	CG-CD2-NE2	-6.14	101.06	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	16	HIS	CG-CD2-NE2	-6.05	101.14	107.20
2	B	280	HIS	CB-CG-ND1	6.02	131.73	122.70
2	D	178	HIS	CB-CG-CD2	-5.91	123.52	131.20
2	D	280	HIS	CB-CG-ND1	5.83	131.45	122.70
2	B	279	VAL	CB-CA-C	-5.80	102.05	110.63
2	B	211	ARG	NE-CZ-NH1	5.79	127.29	121.50
2	B	227	THR	N-CA-CB	-5.77	101.66	110.03
1	C	49	ARG	CA-CB-CG	-5.72	102.65	114.10
1	A	101	HIS	CB-CG-CD2	-5.64	123.87	131.20
2	D	279	VAL	CB-CA-C	-5.64	102.29	110.63
2	D	211	ARG	NE-CZ-NH1	5.62	127.12	121.50
1	C	84	HIS	CB-CG-CD2	-5.59	123.93	131.20
1	A	84	HIS	CB-CG-CD2	-5.58	123.95	131.20
1	C	3	ASN	OD1-CG-ND2	-5.55	117.05	122.60
1	A	53	TYR	O-C-N	-5.53	116.26	122.12
1	C	101	HIS	CB-CG-CD2	-5.53	124.02	131.20
2	B	281	ASP	CA-CB-CG	5.50	118.10	112.60
2	B	262	LYS	N-CA-C	5.47	118.00	111.71
1	A	116	GLN	OE1-CD-NE2	-5.37	117.23	122.60
2	B	205	HIS	CB-CG-CD2	-5.34	124.25	131.20
2	D	161	GLN	N-CA-C	-5.30	106.31	112.89
2	B	161	GLN	N-CA-C	-5.29	106.77	113.18
1	C	74	TYR	N-CA-C	5.23	119.24	112.86
1	C	59	ARG	CB-CA-C	-5.21	101.82	110.68
1	A	16	HIS	CB-CG-CD2	-5.17	124.47	131.20
1	C	56	ARG	CA-C-N	5.15	125.15	120.21
1	C	56	ARG	C-N-CA	5.15	125.15	120.21
1	A	22	TRP	CE2-CD2-CG	-5.13	101.05	107.20
1	C	116	GLN	OE1-CD-NE2	-5.11	117.49	122.60
1	C	22	TRP	CE2-CD2-CG	-5.05	101.14	107.20
1	A	91	VAL	N-CA-C	-5.03	105.64	110.72
1	A	48	GLU	CA-C-O	5.02	126.14	120.92
2	D	276	ASN	CB-CG-ND2	5.01	123.92	116.40
2	B	216	HIS	CB-CG-CD2	-5.00	124.70	131.20

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	1056	0	1052	29	0
1	C	1056	0	1052	25	0
2	B	1037	0	1039	20	0
2	D	1037	0	1039	22	0
3	A	69	0	0	4	0
3	B	52	0	0	4	0
3	C	71	0	0	1	0
3	D	54	0	0	3	0
All	All	4432	0	4182	74	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 9.

All (74) 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:69:ILE:HG23	1:A:96:MET:HE3	1.54	0.89
1:C:82:MET:HE1	1:C:91:VAL:HG11	1.55	0.88
1:C:69:ILE:HG23	1:C:96:MET:HE3	1.54	0.87
1:A:82:MET:HE2	1:A:109:ALA:HB1	1.68	0.74
1:A:101:HIS:HD2	2:D:177:MET:HE2	1.53	0.73
2:B:216:HIS:HD2	2:D:216:HIS:HD2	1.35	0.72
2:B:208:GLU:HG3	2:B:253:ILE:HG23	1.74	0.68
2:B:177:MET:HE3	1:C:104:LEU:HD21	1.79	0.64
1:A:76:ILE:HG13	1:A:102:VAL:HG22	1.81	0.63
1:A:82:MET:HE1	1:A:91:VAL:HG11	1.81	0.63
1:A:76:ILE:HA	1:A:96:MET:HE2	1.82	0.62
3:A:668:HOH:O	2:B:282:GLN:HG2	2.00	0.62
1:C:76:ILE:HA	1:C:96:MET:HE2	1.81	0.61
2:B:189:ALA:O	2:B:215:THR:HB	1.99	0.61
1:C:69:ILE:HG23	1:C:96:MET:CE	2.30	0.61
2:D:189:ALA:O	2:D:215:THR:HB	1.99	0.61
1:A:69:ILE:HG23	1:A:96:MET:CE	2.29	0.59
1:A:68:CYS:HB3	2:B:185:PRO:HA	1.86	0.57
1:C:53:TYR:HB2	1:C:86:LYS:HG3	1.89	0.55
1:A:8:LEU:HD22	2:B:279:VAL:HG13	1.89	0.54
1:A:104:LEU:HD21	2:D:177:MET:HE3	1.90	0.54
1:C:12:ASN:HA	2:D:287:LEU:HD11	1.91	0.53
1:C:76:ILE:HG13	1:C:102:VAL:HG22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:216:HIS:HE1	3:D:683:HOH:O	1.93	0.52
1:C:8:LEU:HD11	2:D:266:TYR:HB2	1.92	0.51
1:C:84:HIS:HB3	1:C:120:LYS:HG3	1.93	0.50
1:A:63:VAL:O	1:A:83:GLU:HG2	2.12	0.50
2:B:249:ASN:HD22	2:B:251:LYS:H	1.58	0.50
1:C:8:LEU:HD22	2:D:279:VAL:HG13	1.93	0.50
1:A:99:THR:HG21	2:D:177:MET:HE1	1.94	0.49
2:B:211:ARG:NH2	3:B:617:HOH:O	2.43	0.49
1:A:101:HIS:CD2	2:D:177:MET:HE2	2.40	0.49
1:A:7:VAL:HG22	1:A:25:LEU:HD12	1.95	0.48
1:A:8:LEU:HD11	2:B:266:TYR:HB2	1.94	0.48
1:C:82:MET:HE2	1:C:109:ALA:C	2.39	0.48
1:C:7:VAL:HG22	1:C:25:LEU:HD12	1.95	0.48
1:C:62:ARG:HD3	1:C:83:GLU:OE2	2.14	0.48
2:D:195:ASN:O	2:D:262:LYS:HD3	2.14	0.47
3:B:653:HOH:O	2:D:216:HIS:HE1	1.97	0.47
1:C:76:ILE:CA	1:C:96:MET:HE2	2.45	0.47
1:A:64:THR:HG22	1:A:83:GLU:HG3	1.97	0.47
1:A:102:VAL:HG13	3:A:717:HOH:O	2.14	0.47
1:A:79:VAL:HB	1:A:82:MET:HE3	1.97	0.47
1:A:86:LYS:NZ	3:A:719:HOH:O	2.48	0.46
1:A:76:ILE:CA	1:A:96:MET:HE2	2.45	0.46
2:B:177:MET:HE2	1:C:101:HIS:ND1	2.31	0.46
1:C:56:ARG:HD2	3:C:677:HOH:O	2.15	0.46
2:D:211:ARG:NH2	3:D:659:HOH:O	2.44	0.45
1:C:68:CYS:HB3	2:D:185:PRO:HA	1.99	0.45
2:B:227:THR:HG23	3:B:722:HOH:O	2.17	0.45
2:B:241:LYS:NZ	3:B:761:HOH:O	2.52	0.43
1:A:85:ILE:HG21	1:A:91:VAL:HG21	2.00	0.43
1:A:104:LEU:HD21	2:D:177:MET:CE	2.49	0.43
1:A:7:VAL:HG13	2:B:157:ALA:HB2	2.01	0.43
1:A:76:ILE:HG13	1:A:102:VAL:CG2	2.47	0.42
2:B:216:HIS:HD2	2:D:216:HIS:CD2	2.24	0.42
1:A:119:LYS:HB3	3:A:649:HOH:O	2.19	0.42
1:C:59:ARG:HD2	2:D:177:MET:HA	2.00	0.42
1:C:7:VAL:HG13	2:D:157:ALA:HB2	2.01	0.42
1:A:80:ALA:HB3	1:A:105:VAL:HG12	2.02	0.41
2:B:162:GLY:O	2:B:262:LYS:HE3	2.20	0.41
2:B:227:THR:HG22	2:B:230:GLN:H	1.84	0.41
2:D:271:ILE:HG22	2:D:272:GLN:HG3	2.02	0.41
1:A:82:MET:CE	1:A:109:ALA:HB1	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:216:HIS:CD2	2:D:216:HIS:HD2	2.25	0.41
1:C:59:ARG:HG3	2:D:176:LYS:O	2.20	0.41
1:C:82:MET:CE	1:C:109:ALA:HB1	2.51	0.41
2:D:205:HIS:HB3	3:D:765:HOH:O	2.20	0.41
1:C:80:ALA:HB3	1:C:105:VAL:HG12	2.02	0.41
1:A:72:GLU:O	2:B:195:ASN:HB2	2.21	0.40
1:A:130:LYS:O	1:A:134:GLU:HG3	2.21	0.40
1:C:49:ARG:HH11	1:C:49:ARG:HD3	1.67	0.40
1:C:130:LYS:O	1:C:134:GLU:HG3	2.21	0.40
2:D:187:ILE:O	2:D:191:LEU:O	2.38	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	134/151 (89%)	131 (98%)	3 (2%)	0	100	100
1	C	134/151 (89%)	131 (98%)	3 (2%)	0	100	100
2	B	137/144 (95%)	133 (97%)	4 (3%)	0	100	100
2	D	137/144 (95%)	132 (96%)	5 (4%)	0	100	100
All	All	542/590 (92%)	527 (97%)	15 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	113/128 (88%)	106 (94%)	7 (6%)	16	14
1	C	113/128 (88%)	106 (94%)	7 (6%)	16	14
2	B	107/110 (97%)	100 (94%)	7 (6%)	15	13
2	D	107/110 (97%)	100 (94%)	7 (6%)	15	13
All	All	440/476 (92%)	412 (94%)	28 (6%)	16	14

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	8	LEU
1	A	16	HIS
1	A	102	VAL
1	A	119	LYS
1	A	121	GLU
1	A	138	THR
2	B	185	PRO
2	B	201	THR
2	B	215	THR
2	B	227	THR
2	B	234	GLU
2	B	249	ASN
2	B	254	GLN
1	C	7	VAL
1	C	8	LEU
1	C	16	HIS
1	C	102	VAL
1	C	119	LYS
1	C	123	LEU
1	C	138	THR
2	D	185	PRO
2	D	201	THR
2	D	215	THR
2	D	234	GLU
2	D	243	VAL
2	D	254	GLN
2	D	262	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	ASN
2	B	216	HIS
2	B	223	ASN
2	B	249	ASN
2	B	276	ASN
1	C	12	ASN
1	C	16	HIS
1	C	73	ASN
1	C	87	ASN
2	D	161	GLN
2	D	216	HIS
2	D	223	ASN
2	D	276	ASN

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	136/151 (90%)	-0.17	1 (0%) 84 86	8, 14, 29, 35	0
1	C	136/151 (90%)	-0.14	0 100 100	8, 14, 29, 35	0
2	B	139/144 (96%)	-0.25	0 100 100	7, 13, 20, 22	0
2	D	139/144 (96%)	-0.20	2 (1%) 73 75	7, 14, 20, 23	0
All	All	550/590 (93%)	-0.19	3 (0%) 87 88	7, 14, 24, 35	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	172	GLY	2.9
1	A	16	HIS	2.6
2	D	273	ASP	2.5

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.