



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 03:16 AM EDT

PDB ID : 1WUV / pdb_00001wuv
Title : Crystal structure of native Canavalia gladiata lectin (CGL): a tetrameric ConA-like lectin
Authors : Freitas, B.T.; Delatorre, P.; Moreno, F.B.M.B.; Rocha, B.A.M.; Souza, E.P.; Canduri, F.; Cardoso, A.L.H.; Sampaio, A.H.; Azevedo Jr., W.F.; Cavada, B.S.
Deposited on : 2004-12-09
Resolution : 2.30 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

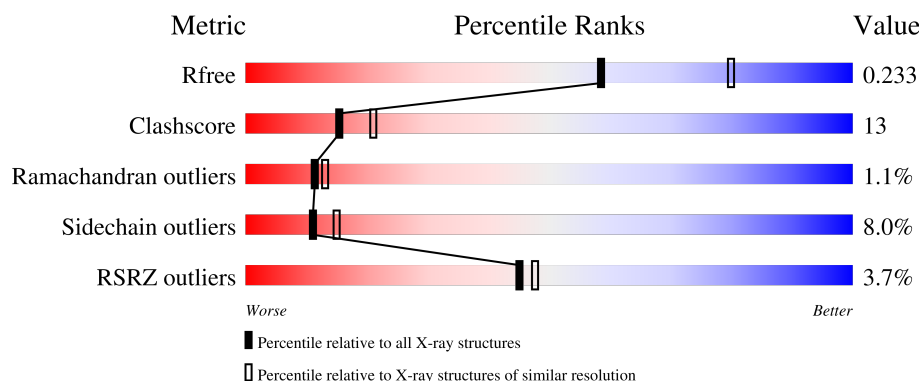
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	237	5% 75% 19% 5%
1	D	237	4% 64% 30%
1	G	237	3% 68% 27% 5%
1	J	237	3% 68% 25% 5%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7640 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 Concanavalin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1805	1138	303	362	2			
1	D	237	Total	C	N	O	S	0	0	0
			1805	1138	303	362	2			
1	G	237	Total	C	N	O	S	0	0	0
			1805	1138	303	362	2			
1	J	237	Total	C	N	O	S	0	0	0
			1805	1138	303	362	2			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	D	1	Total	Mn	0	0
			1	1		
2	G	1	Total	Mn	0	0
			1	1		
2	J	1	Total	Mn	0	0
			1	1		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		
3	G	1	Total	Ca	0	0
			1	1		
3	J	1	Total	Ca	0	0
			1	1		

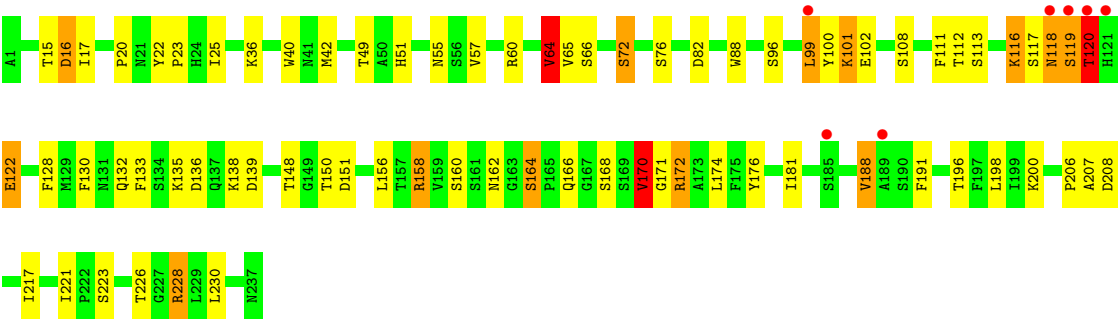
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	114	Total 114	O 114	0	0
4	D	101	Total 101	O 101	0	0
4	G	99	Total 99	O 99	0	0
4	J	98	Total 98	O 98	0	0

- Molecule 1: Concanavalin A



● Molecule 1: Concanavalin A



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	100.90Å 115.35Å 241.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.30 10.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (10.00-2.30) 96.5 (10.00-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.184 , 0.233 0.187 , 0.233	Depositor DCC
R_{free} test set	3062 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7640	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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: CA, MN

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	1.24	3/1847 (0.2%)	1.25	8/2516 (0.3%)
1	D	1.25	3/1847 (0.2%)	1.29	12/2516 (0.5%)
1	G	1.20	7/1847 (0.4%)	1.15	2/2516 (0.1%)
1	J	1.22	2/1847 (0.1%)	1.17	8/2516 (0.3%)
All	All	1.23	15/7388 (0.2%)	1.21	30/10064 (0.3%)

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	D	0	2
1	G	0	1
1	J	0	1
All	All	0	4

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	16	ASP	CB-CG	7.08	1.69	1.52
1	G	120	THR	CA-CB	6.81	1.65	1.53
1	J	120	THR	CA-CB	6.12	1.63	1.53
1	A	189	ALA	C-O	5.89	1.30	1.23
1	A	120	THR	CA-CB	5.74	1.63	1.53
1	G	43	GLN	CG-CD	5.61	1.66	1.52
1	G	54	TYR	C-O	-5.56	1.17	1.23
1	D	172	ARG	CD-NE	-5.38	1.38	1.46
1	G	188	VAL	CA-CB	5.33	1.61	1.54
1	J	99	LEU	CB-CG	5.21	1.63	1.53
1	G	120	THR	N-CA	5.14	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	194	THR	C-O	-5.12	1.17	1.23
1	D	7	VAL	CA-CB	-5.08	1.48	1.54
1	G	89	VAL	CA-CB	5.08	1.62	1.54
1	A	172	ARG	CD-NE	-5.04	1.39	1.46

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	203	ASP	CA-C-N	-9.43	109.30	122.87
1	D	203	ASP	C-N-CA	-9.43	109.30	122.87
1	J	64	VAL	CB-CA-C	-8.93	96.75	110.50
1	D	64	VAL	CB-CA-C	-8.78	96.98	110.50
1	G	64	VAL	CB-CA-C	-8.53	97.36	110.50
1	D	119	SER	N-CA-C	8.29	121.10	110.65
1	A	64	VAL	CB-CA-C	-7.77	98.54	110.50
1	J	119	SER	N-CA-C	7.69	122.08	111.55
1	A	57	VAL	N-CA-C	-7.36	104.61	111.45
1	D	189	ALA	N-CA-C	-7.16	98.18	109.50
1	D	217	ILE	N-CA-C	6.48	117.23	110.62
1	A	170	VAL	CB-CA-C	-6.43	99.58	110.71
1	D	203	ASP	N-CA-C	-6.25	97.48	110.80
1	J	101	LYS	N-CA-C	6.06	118.16	109.14
1	A	22	TYR	N-CA-C	-5.67	103.03	108.13
1	D	170	VAL	CB-CA-C	-5.57	100.74	110.71
1	A	65	VAL	CB-CA-C	-5.56	102.33	110.62
1	D	172	ARG	CB-CG-CD	-5.54	98.57	111.30
1	D	159	VAL	N-CA-C	-5.51	100.18	108.12
1	J	16	ASP	N-CA-CB	-5.47	101.24	110.49
1	G	119	SER	N-CA-C	5.42	122.35	110.80
1	A	150	THR	OG1-CB-CG2	-5.36	98.57	109.30
1	J	170	VAL	CB-CA-C	-5.34	102.28	110.50
1	A	123	THR	CB-CA-C	-5.25	102.60	110.16
1	D	214	ILE	N-CA-C	-5.25	100.50	108.17
1	J	15	THR	CA-C-N	5.21	131.49	121.54
1	J	15	THR	C-N-CA	5.21	131.49	121.54
1	A	170	VAL	CA-CB-CG2	5.20	119.23	110.40
1	J	72	SER	N-CA-C	5.04	116.89	108.99
1	D	228	ARG	NE-CZ-NH2	-5.00	114.70	119.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	118	ASN	Peptide
1	D	123	THR	Peptide
1	G	118	ASN	Peptide
1	J	118	ASN	Peptide

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	1805	0	1751	37	0
1	D	1805	0	1751	54	0
1	G	1805	0	1751	56	0
1	J	1805	0	1751	47	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
2	G	1	0	0	0	0
2	J	1	0	0	0	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
3	G	1	0	0	0	0
3	J	1	0	0	0	0
4	A	114	0	0	0	0
4	D	101	0	0	1	0
4	G	99	0	0	1	0
4	J	98	0	0	1	0
All	All	7640	0	7004	192	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 13.

All (192) 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:G:43:GLN:NE2	1:G:69:ASN:HD21	1.16	1.41
1:G:43:GLN:HE22	1:G:69:ASN:ND2	1.24	1.32
1:J:119:SER:O	1:J:122:GLU:HG2	1.51	1.07
1:G:42:MET:HE1	1:G:206:PRO:HG2	1.37	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:25:ILE:HD12	1:G:65:VAL:HG23	1.46	0.97
1:D:160:SER:HB3	1:D:166:GLN:NE2	1.81	0.95
1:D:25:ILE:HD12	1:D:65:VAL:HG23	1.49	0.92
1:G:42:MET:HE1	1:G:206:PRO:CG	2.01	0.91
1:J:170:VAL:HG22	1:J:226:THR:HG22	1.55	0.88
1:J:25:ILE:HD12	1:J:65:VAL:HG23	1.56	0.87
1:J:160:SER:HB3	1:J:166:GLN:NE2	1.89	0.86
1:G:43:GLN:NE2	1:G:69:ASN:ND2	1.97	0.85
1:A:25:ILE:HD12	1:A:65:VAL:HG23	1.58	0.84
1:A:119:SER:HB2	1:A:122:GLU:HG3	1.60	0.83
1:J:119:SER:C	1:J:122:GLU:HG2	2.02	0.83
1:G:25:ILE:HD12	1:G:65:VAL:CG2	2.09	0.83
1:D:160:SER:HB3	1:D:166:GLN:HE22	1.41	0.82
1:G:42:MET:CE	1:G:206:PRO:HG2	2.10	0.80
1:J:119:SER:HB2	1:J:122:GLU:HG3	1.62	0.80
1:A:25:ILE:HD12	1:A:65:VAL:CG2	2.13	0.79
1:D:25:ILE:HD12	1:D:65:VAL:CG2	2.12	0.79
1:A:170:VAL:HG22	1:A:226:THR:HG22	1.63	0.79
1:J:88:TRP:HB3	1:J:217:ILE:HD11	1.66	0.78
1:G:119:SER:HB2	1:G:122:GLU:HG3	1.66	0.76
1:D:156:LEU:O	1:D:171:GLY:HA3	1.87	0.75
1:D:202:PRO:C	1:D:203:ASP:O	2.24	0.75
1:G:66:SER:HB3	1:G:72:SER:CB	2.17	0.75
1:G:172:ARG:HD2	1:G:221:ILE:HG12	1.69	0.74
1:J:16:ASP:OD2	1:J:228:ARG:NH2	2.24	0.70
1:A:87:GLU:HG2	1:A:182:TRP:O	1.91	0.69
1:D:66:SER:HB3	1:D:72:SER:CB	2.22	0.69
1:D:170:VAL:HG22	1:D:226:THR:HG22	1.74	0.68
1:D:88:TRP:HB3	1:D:217:ILE:HD11	1.77	0.67
1:G:116:LYS:HD3	1:G:123:THR:OG1	1.94	0.67
1:G:162:ASN:OD1	1:G:164:SER:HB2	1.95	0.67
1:A:137:GLN:HG2	1:A:140:LEU:HD12	1.76	0.66
1:J:66:SER:HB3	1:J:72:SER:HB2	1.76	0.66
1:D:87:GLU:HG3	1:D:182:TRP:O	1.96	0.66
1:D:66:SER:HB3	1:D:72:SER:HB3	1.77	0.66
1:J:42:MET:HE1	1:J:206:PRO:HG3	1.76	0.66
1:G:66:SER:HB3	1:G:72:SER:HB2	1.78	0.66
1:G:118:ASN:ND2	1:G:118:ASN:H	1.94	0.65
1:D:228:ARG:HD2	4:D:268:HOH:O	1.96	0.65
1:J:160:SER:HB3	1:J:166:GLN:HE21	1.60	0.64
1:G:118:ASN:H	1:G:118:ASN:HD22	1.43	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:116:LYS:HG3	1:G:188:VAL:HB	1.80	0.64
1:A:80:ASP:OD1	1:A:82:ASP:HB2	2.00	0.62
1:D:170:VAL:HG22	1:D:226:THR:CG2	2.30	0.62
1:J:42:MET:CE	1:J:206:PRO:HG3	2.30	0.62
1:J:17:ILE:HD13	1:J:228:ARG:HD3	1.83	0.61
1:G:66:SER:HB3	1:G:72:SER:HB3	1.84	0.59
1:J:66:SER:HB3	1:J:72:SER:CB	2.32	0.59
1:A:160:SER:HB3	1:A:166:GLN:NE2	2.18	0.59
1:J:158:ARG:HH11	1:J:158:ARG:HG3	1.68	0.59
1:J:60:ARG:HD3	1:J:76:SER:HB3	1.85	0.59
1:D:118:ASN:ND2	1:D:118:ASN:H	2.00	0.58
1:J:172:ARG:HD2	1:J:221:ILE:HG12	1.84	0.58
1:G:120:THR:HG22	1:G:121:HIS:HD2	1.67	0.58
1:A:170:VAL:HG22	1:A:226:THR:CG2	2.33	0.57
1:D:137:GLN:NE2	1:D:139:ASP:OD1	2.28	0.57
1:G:119:SER:O	1:G:122:GLU:HG2	2.05	0.56
1:D:133:PHE:O	1:D:152:GLY:HA2	2.06	0.56
1:D:25:ILE:CD1	1:D:65:VAL:HG23	2.29	0.56
1:G:30:LYS:HD2	1:G:84:VAL:HG13	1.88	0.56
1:D:66:SER:HB3	1:D:72:SER:HB2	1.88	0.55
1:A:20:PRO:HB2	1:A:22:TYR:CZ	2.42	0.54
1:G:43:GLN:HE22	1:G:69:ASN:HD21	0.56	0.54
1:A:90:ARG:NH1	1:A:217:ILE:HG23	2.23	0.54
1:J:160:SER:HB3	1:J:166:GLN:HE22	1.72	0.54
1:A:42:MET:HE1	1:A:206:PRO:HG2	1.89	0.53
1:J:172:ARG:HD2	1:J:221:ILE:CG1	2.39	0.53
1:J:162:ASN:OD1	1:J:164:SER:HB3	2.09	0.53
1:J:49:THR:OG1	1:J:196:THR:HG22	2.09	0.53
1:A:172:ARG:HD2	1:A:221:ILE:CG1	2.39	0.52
1:J:119:SER:HB2	1:J:122:GLU:CG	2.35	0.52
1:G:55:ASN:HB3	4:G:256:HOH:O	2.09	0.52
1:G:17:ILE:HD13	1:G:228:ARG:HD3	1.91	0.52
1:A:88:TRP:HB3	1:A:217:ILE:HD11	1.92	0.51
1:G:88:TRP:HB3	1:G:217:ILE:HD11	1.91	0.51
1:D:172:ARG:HD2	1:D:221:ILE:CG1	2.41	0.51
1:J:116:LYS:HG3	1:J:188:VAL:HB	1.92	0.51
1:A:118:ASN:HA	1:A:187:VAL:HG21	1.93	0.51
1:A:172:ARG:HD2	1:A:221:ILE:HG13	1.93	0.51
1:A:160:SER:HB3	1:A:166:GLN:HE21	1.76	0.51
1:J:23:PRO:HB2	1:J:40:TRP:O	2.10	0.51
1:A:96:SER:OG	1:A:230:LEU:HA	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:183:GLU:HB3	1:D:186:ALA:HB2	1.94	0.49
1:J:51:HIS:HB2	1:J:64:VAL:HG23	1.95	0.49
1:A:56:SER:HB3	1:A:189:ALA:O	2.13	0.49
1:A:119:SER:HB2	1:A:122:GLU:CG	2.38	0.49
1:G:43:GLN:HE22	1:G:69:ASN:CG	2.09	0.49
1:J:100:TYR:HB2	1:J:207:ALA:CB	2.43	0.49
1:D:160:SER:HB3	1:D:166:GLN:HE21	1.73	0.49
1:D:51:HIS:HB2	1:D:64:VAL:HG23	1.95	0.48
1:G:87:GLU:HG3	1:G:182:TRP:O	2.13	0.48
1:G:42:MET:HE1	1:G:206:PRO:HG3	1.92	0.48
1:G:45:GLY:HA2	1:G:200:LYS:HG3	1.95	0.48
1:G:96:SER:OG	1:G:230:LEU:HA	2.13	0.48
1:J:158:ARG:HG3	1:J:158:ARG:NH1	2.29	0.48
1:G:116:LYS:HD3	1:G:123:THR:HG1	1.76	0.48
1:G:156:LEU:O	1:G:171:GLY:HA3	2.13	0.48
1:D:4:ILE:HD13	1:D:215:SER:HB3	1.95	0.47
1:D:14:ASN:O	1:D:19:ASP:HB2	2.14	0.47
1:G:118:ASN:ND2	1:G:185:SER:O	2.47	0.47
1:J:172:ARG:NH2	4:J:251:HOH:O	2.45	0.47
1:A:42:MET:HB3	1:A:42:MET:HE3	1.50	0.47
1:D:122:GLU:HA	1:D:122:GLU:OE1	2.15	0.47
1:D:172:ARG:HD2	1:D:221:ILE:HG13	1.96	0.47
1:G:170:VAL:HG22	1:G:226:THR:HG22	1.97	0.47
1:A:208:ASP:OD1	1:A:227:GLY:HA2	2.15	0.47
1:G:90:ARG:NH1	1:G:217:ILE:HG23	2.30	0.47
1:D:49:THR:OG1	1:D:196:THR:HG22	2.15	0.47
1:G:14:ASN:O	1:G:19:ASP:HB2	2.14	0.47
1:G:91:VAL:HG12	1:G:214:ILE:HG12	1.95	0.47
1:D:42:MET:CE	1:D:206:PRO:HG3	2.44	0.47
1:D:111:PHE:CE2	1:D:113:SER:HB2	2.50	0.47
1:D:51:HIS:O	1:D:63:ALA:HA	2.14	0.47
1:D:20:PRO:HB2	1:D:22:TYR:CZ	2.50	0.47
1:J:102:GLU:HG2	1:J:207:ALA:O	2.15	0.47
1:J:136:ASP:OD2	1:J:138:LYS:NZ	2.43	0.46
1:A:156:LEU:O	1:A:171:GLY:HA3	2.15	0.46
1:D:201:SER:O	1:D:203:ASP:O	2.34	0.46
1:D:122:GLU:OE1	1:D:122:GLU:CA	2.63	0.46
1:J:96:SER:OG	1:J:230:LEU:HA	2.16	0.46
1:A:137:GLN:HG2	1:A:140:LEU:CD1	2.45	0.46
1:G:52:ILE:O	1:G:192:ASP:HA	2.15	0.46
1:D:3:THR:O	1:D:215:SER:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:14:ASN:C	1:G:19:ASP:HB2	2.41	0.46
1:J:162:ASN:OD1	1:J:162:ASN:C	2.59	0.46
1:G:3:THR:O	1:G:215:SER:HA	2.17	0.45
1:A:11:THR:O	1:A:205:HIS:HE1	2.00	0.45
1:D:176:TYR:CE1	1:G:176:TYR:CE1	3.04	0.45
1:J:112:THR:O	1:J:191:PHE:HA	2.17	0.45
1:D:211:ALA:CB	1:D:232:LEU:HD11	2.47	0.44
1:A:133:PHE:HA	1:A:137:GLN:OE1	2.18	0.44
1:D:126:LEU:CD2	1:D:179:VAL:HG22	2.48	0.44
1:D:119:SER:HB2	1:D:122:GLU:HG2	2.00	0.44
1:J:100:TYR:HB2	1:J:207:ALA:HB3	2.00	0.43
1:G:20:PRO:HB2	1:G:22:TYR:CE1	2.53	0.43
1:J:25:ILE:HD12	1:J:65:VAL:CG2	2.36	0.43
1:A:57:VAL:HG23	1:A:188:VAL:HG13	2.00	0.43
1:A:68:PRO:O	1:A:69:ASN:C	2.61	0.43
1:A:106:ILE:HB	1:A:154:LEU:HB3	2.00	0.43
1:D:126:LEU:HD23	1:D:179:VAL:HG22	2.01	0.43
1:D:105:THR:O	1:D:197:PHE:HA	2.18	0.43
1:G:9:LEU:HD11	1:G:65:VAL:HG22	2.01	0.43
1:G:92:GLY:HA2	1:G:109:TRP:CH2	2.54	0.43
1:G:172:ARG:HD2	1:G:221:ILE:CG1	2.44	0.43
1:J:139:ASP:O	1:J:176:TYR:HB2	2.19	0.43
1:D:43:GLN:HE21	1:D:46:LYS:HG3	1.84	0.42
1:D:187:VAL:HG12	1:D:188:VAL:HG23	2.00	0.42
1:A:133:PHE:HB3	1:A:148:THR:HB	2.01	0.42
1:J:111:PHE:CE2	1:J:113:SER:HB2	2.55	0.42
1:J:117:SER:O	1:J:119:SER:N	2.52	0.42
1:J:108:SER:HA	1:J:130:PHE:O	2.18	0.42
1:J:111:PHE:HB3	1:J:128:PHE:CZ	2.55	0.42
1:J:174:LEU:HD12	1:J:174:LEU:N	2.35	0.42
1:D:124:ASN:HA	1:G:129:MET:O	2.20	0.42
1:D:117:SER:HA	1:D:186:ALA:HA	2.02	0.42
1:J:133:PHE:HB3	1:J:148:THR:HB	2.02	0.42
1:D:14:ASN:C	1:D:19:ASP:HB2	2.45	0.42
1:G:20:PRO:HB2	1:G:22:TYR:CZ	2.55	0.42
1:A:207:ALA:HA	1:A:208:ASP:HA	1.94	0.41
1:G:211:ALA:CB	1:G:232:LEU:HD11	2.50	0.41
1:J:20:PRO:HB2	1:J:22:TYR:CZ	2.55	0.41
1:D:133:PHE:HB3	1:D:148:THR:HB	2.00	0.41
1:G:102:GLU:HB2	1:G:199:ILE:HG23	2.02	0.41
1:A:23:PRO:HB2	1:A:40:TRP:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ILE:HD13	1:A:228:ARG:HD3	2.02	0.41
1:A:170:VAL:HG11	1:A:231:GLY:HA2	2.03	0.41
1:D:12:TYR:HA	1:D:13:PRO:HD3	1.82	0.41
1:A:3:THR:O	1:A:215:SER:HA	2.21	0.41
1:D:68:PRO:O	1:D:69:ASN:C	2.64	0.41
1:D:208:ASP:OD1	1:D:227:GLY:HA2	2.20	0.41
1:G:111:PHE:CE2	1:G:113:SER:HB2	2.55	0.41
1:J:42:MET:HE3	1:J:42:MET:HB3	1.89	0.41
1:A:205:HIS:ND1	1:A:206:PRO:HD2	2.36	0.41
1:D:68:PRO:C	1:D:70:GLY:N	2.79	0.41
1:G:52:ILE:C	1:G:53:ILE:HG13	2.46	0.41
1:D:52:ILE:O	1:D:192:ASP:HA	2.21	0.41
1:D:135:LYS:HB3	1:D:135:LYS:HE2	1.70	0.41
1:G:118:ASN:ND2	1:G:118:ASN:N	2.63	0.41
1:J:156:LEU:O	1:J:171:GLY:HA3	2.21	0.41
1:G:11:THR:HB	1:G:42:MET:HE2	2.03	0.40
1:G:207:ALA:HA	1:G:208:ASP:HA	1.86	0.40
1:D:181:ILE:HD12	1:D:181:ILE:HA	1.76	0.40
1:J:207:ALA:HA	1:J:208:ASP:HA	1.80	0.40
1:A:55:ASN:OD1	1:A:57:VAL:N	2.49	0.40
1:D:91:VAL:HG12	1:D:214:ILE:HG12	2.03	0.40
1:G:170:VAL:HG13	1:G:226:THR:HA	2.04	0.40
1:J:55:ASN:OD1	1:J:57:VAL:N	2.53	0.40
1:G:28:ASP:HB3	1:G:31:SER:O	2.21	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	235/237 (99%)	219 (93%)	13 (6%)	3 (1%)	9 10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	235/237 (99%)	219 (93%)	13 (6%)	3 (1%)	9	10
1	G	235/237 (99%)	225 (96%)	8 (3%)	2 (1%)	14	17
1	J	235/237 (99%)	222 (94%)	11 (5%)	2 (1%)	14	17
All	All	940/948 (99%)	885 (94%)	45 (5%)	10 (1%)	11	13

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	D	120	THR
1	D	203	ASP
1	G	120	THR
1	D	150	THR
1	J	118	ASN
1	A	120	THR
1	A	122	GLU
1	J	120	THR
1	G	118	ASN

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	202/202 (100%)	190 (94%)	12 (6%)	18	26
1	D	202/202 (100%)	186 (92%)	16 (8%)	11	15
1	G	202/202 (100%)	188 (93%)	14 (7%)	14	20
1	J	202/202 (100%)	179 (89%)	23 (11%)	5	6
All	All	808/808 (100%)	743 (92%)	65 (8%)	11	15

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ASN

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Mol	Chain	Res	Type
1	A	64	VAL
1	A	69	ASN
1	A	71	ASP
1	A	99	LEU
1	A	118	ASN
1	A	119	SER
1	A	123	THR
1	A	170	VAL
1	A	181	ILE
1	A	188	VAL
1	A	200	LYS
1	D	64	VAL
1	D	101	LYS
1	D	116	LYS
1	D	118	ASN
1	D	121	HIS
1	D	122	GLU
1	D	150	THR
1	D	161	SER
1	D	164	SER
1	D	170	VAL
1	D	181	ILE
1	D	187	VAL
1	D	188	VAL
1	D	198	LEU
1	D	200	LYS
1	D	228	ARG
1	G	9	LEU
1	G	39	LYS
1	G	64	VAL
1	G	66	SER
1	G	99	LEU
1	G	116	LYS
1	G	132	GLN
1	G	135	LYS
1	G	161	SER
1	G	164	SER
1	G	170	VAL
1	G	187	VAL
1	G	200	LYS
1	G	228	ARG
1	J	36	LYS

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Mol	Chain	Res	Type
1	J	64	VAL
1	J	82	ASP
1	J	99	LEU
1	J	101	LYS
1	J	116	LYS
1	J	120	THR
1	J	122	GLU
1	J	132	GLN
1	J	135	LYS
1	J	150	THR
1	J	151	ASP
1	J	158	ARG
1	J	164	SER
1	J	168	SER
1	J	170	VAL
1	J	172	ARG
1	J	181	ILE
1	J	188	VAL
1	J	198	LEU
1	J	200	LYS
1	J	223	SER
1	J	228	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	132	GLN
1	D	21	ASN
1	D	43	GLN
1	D	118	ASN
1	D	132	GLN
1	D	166	GLN
1	G	21	ASN
1	G	43	GLN
1	G	118	ASN
1	G	121	HIS
1	G	132	GLN
1	J	43	GLN
1	J	118	ASN
1	J	121	HIS

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](#)

Of 8 ligands modelled in this entry, 8 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 [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	237/237 (100%)	-0.48	11 (4%) 37 39	8, 15, 39, 63	0
1	D	237/237 (100%)	-0.50	10 (4%) 40 42	8, 15, 40, 59	0
1	G	237/237 (100%)	-0.39	7 (2%) 52 54	10, 20, 41, 58	0
1	J	237/237 (100%)	-0.35	7 (2%) 52 54	8, 18, 41, 57	0
All	All	948/948 (100%)	-0.43	35 (3%) 45 48	8, 18, 41, 63	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	118	ASN	5.7
1	J	118	ASN	5.4
1	A	119	SER	5.4
1	A	118	ASN	5.3
1	D	118	ASN	5.2
1	G	119	SER	5.1
1	D	120	THR	5.0
1	J	120	THR	4.8
1	J	185	SER	4.7
1	D	16	ASP	4.4
1	G	185	SER	4.4
1	G	120	THR	4.3
1	D	119	SER	4.3
1	G	69	ASN	4.1
1	A	120	THR	4.1
1	G	121	HIS	4.1
1	D	121	HIS	4.0
1	A	189	ALA	3.9
1	J	121	HIS	3.8
1	J	119	SER	3.8
1	A	69	ASN	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	184	SER	3.5
1	D	185	SER	3.4
1	A	185	SER	3.3
1	J	99	LEU	3.1
1	D	203	ASP	3.1
1	A	121	HIS	3.0
1	D	69	ASN	3.0
1	D	117	SER	2.7
1	A	122	GLU	2.4
1	A	117	SER	2.3
1	A	82	ASP	2.3
1	D	122	GLU	2.1
1	J	189	ALA	2.0
1	G	43	GLN	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
3	CA	J	240	1/1	0.98	0.02	18,18,18,18	0
3	CA	A	240	1/1	0.99	0.01	11,11,11,11	0
3	CA	D	240	1/1	0.99	0.01	11,11,11,11	0
3	CA	G	240	1/1	0.99	0.01	17,17,17,17	0
2	MN	J	239	1/1	0.99	0.02	17,17,17,17	0
2	MN	G	239	1/1	1.00	0.02	18,18,18,18	0
2	MN	A	239	1/1	1.00	0.03	13,13,13,13	0
2	MN	D	239	1/1	1.00	0.02	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.