



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:26 AM GMT

PDB ID : 2D3R
Title : Cratylia folibunda seed lectin at acidic pH
Authors : Del Sol, F.G.; Cavada, B.S.; Calvete, J.J.
Deposited on : 2005-09-30
Resolution : 2.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

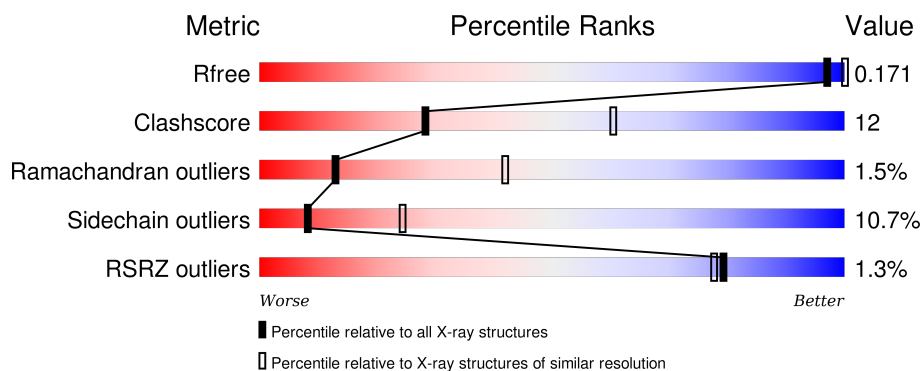
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	236	<div> <div>72%</div> <div>24%</div> <div>.</div> </div>
1	B	236	<div> <div>3%</div> <div>67%</div> <div>28%</div> <div>..</div> </div>
1	C	236	<div> <div>%</div> <div>70%</div> <div>23%</div> <div>6%</div> </div>
1	D	236	<div> <div>%</div> <div>70%</div> <div>25%</div> <div>5%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7258 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 Lectin alpha chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	236	Total	C	N	O	0	0	0
			1794	1126	305	363			
1	B	236	Total	C	N	O	0	0	0
			1794	1126	305	363			
1	C	236	Total	C	N	O	0	0	0
			1794	1126	305	363			
1	D	236	Total	C	N	O	0	0	0
			1794	1126	305	363			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	PHE	TRP	SEE REMARK 999	UNP P81517
B	211	PHE	TRP	SEE REMARK 999	UNP P81517
C	211	PHE	TRP	SEE REMARK 999	UNP P81517
D	211	PHE	TRP	SEE REMARK 999	UNP P81517

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Mn 1	0	0
3	A	1	Total 1	Mn 1	0	0
3	D	1	Total 1	Mn 1	0	0
3	C	1	Total 1	Mn 1	0	0

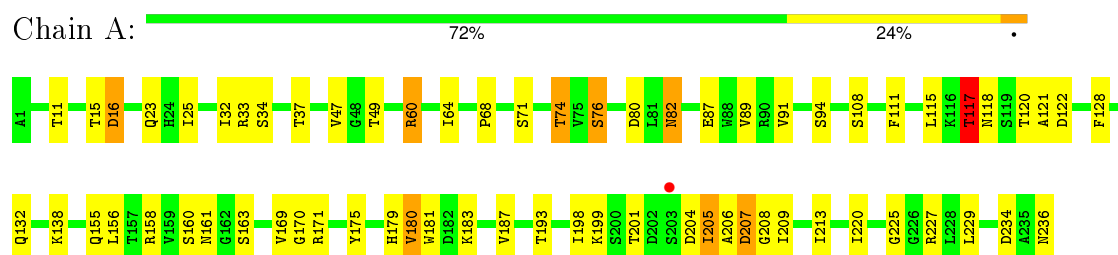
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total 22	O 22	0	0
4	B	21	Total 21	O 21	0	0
4	C	18	Total 18	O 18	0	0
4	D	13	Total 13	O 13	0	0

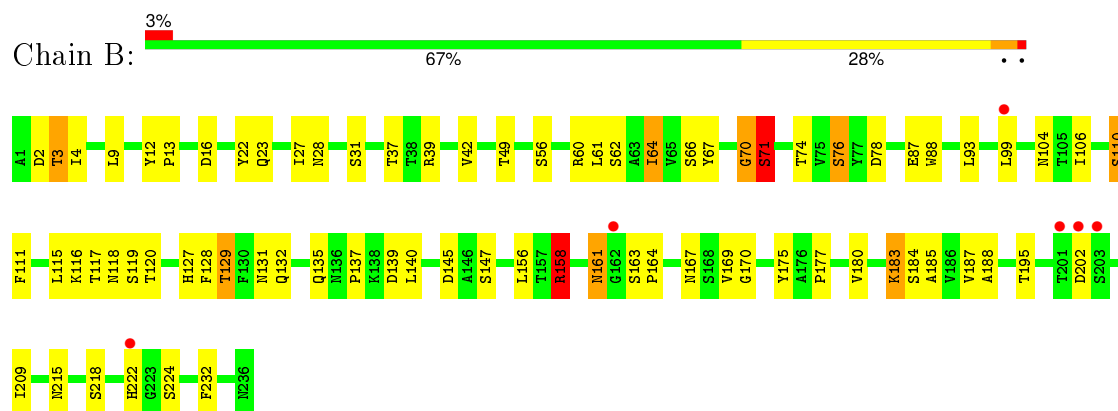
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

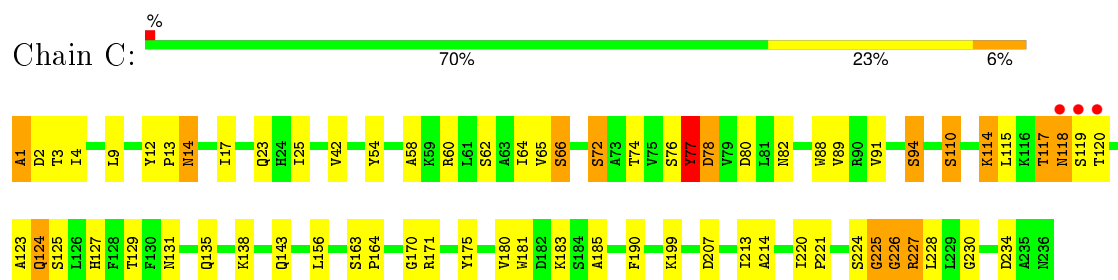
• Molecule 1: Lectin alpha chain



• Molecule 1: Lectin alpha chain

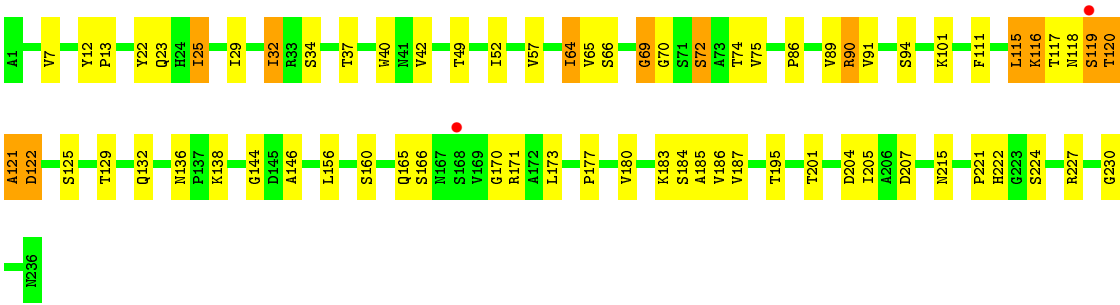


• Molecule 1: Lectin alpha chain



• Molecule 1: Lectin alpha chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.21Å 125.44Å 126.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.80 – 2.90 49.86 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (49.80-2.90) 68.3 (49.86-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.77 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.178 , 0.220 0.184 , 0.171	Depositor DCC
R_{free} test set	1393 reflections (9.79%)	DCC
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 33.0	EDS
Estimated twinning fraction	0.036 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 16570 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7258	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.13% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.52	0/1833	0.70	2/2500 (0.1%)
1	B	0.53	0/1833	0.66	1/2500 (0.0%)
1	C	0.64	1/1833 (0.1%)	0.88	9/2500 (0.4%)
1	D	0.53	0/1833	0.73	1/2500 (0.0%)
All	All	0.56	1/7332 (0.0%)	0.75	13/10000 (0.1%)

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	A	1	0
1	C	0	2
All	All	1	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	227	ARG	CB-CG	7.38	1.72	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	2	ASP	CA-C-N	-13.64	87.20	117.20
1	D	121	ALA	N-CA-C	-10.89	81.58	111.00
1	C	2	ASP	O-C-N	9.32	137.62	122.70
1	C	1	ALA	C-N-CA	8.92	144.01	121.70
1	C	2	ASP	C-N-CA	7.96	141.61	121.70
1	C	2	ASP	CA-C-O	7.18	135.17	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	2	ASP	CA-CB-CG	-7.03	97.94	113.40
1	C	227	ARG	CA-CB-CG	-5.98	100.25	113.40
1	A	117	THR	OG1-CB-CG2	5.64	122.97	110.00
1	A	117	THR	CA-CB-OG1	5.43	120.40	109.00
1	C	78	ASP	N-CA-C	-5.39	96.45	111.00
1	B	158	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	C	77	TYR	C-N-CA	-5.12	108.89	121.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	117	THR	CB

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	77	TYR	Mainchain,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	1794	0	1736	41	0
1	B	1794	0	1736	49	12
1	C	1794	0	1735	51	12
1	D	1794	0	1736	47	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	22	0	0	0	0
4	B	21	0	0	1	0
4	C	18	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	13	0	0	0	0
All	All	7258	0	6943	164	12

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 12.

All (164) 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:D:117:THR:CG2	1:D:185:ALA:HA	1.67	1.23
1:D:117:THR:HG22	1:D:185:ALA:CA	1.75	1.14
1:D:117:THR:HB	1:D:184:SER:O	1.51	1.11
1:C:3:THR:O	1:C:4:ILE:HG12	1.49	1.10
1:C:110:SER:HB3	1:C:129:THR:HG22	1.03	1.02
1:D:144:GLY:HA3	1:D:170:GLY:HA2	1.39	1.00
1:B:195:THR:HG23	1:C:120:THR:HG21	1.43	1.00
1:B:158:ARG:HG3	1:B:158:ARG:HH11	1.24	0.99
1:D:117:THR:HG22	1:D:185:ALA:HA	1.00	0.98
1:B:195:THR:CG2	1:C:120:THR:HG21	1.96	0.96
1:B:60:ARG:NH1	1:C:58:ALA:O	2.03	0.90
1:B:195:THR:CG2	1:C:120:THR:CG2	2.49	0.90
1:B:3:THR:HG23	1:B:215:ASN:HD21	1.41	0.84
1:A:115:LEU:HD21	1:A:179:HIS:O	1.77	0.84
1:B:3:THR:HG23	1:B:215:ASN:ND2	1.94	0.83
1:B:195:THR:HG22	1:C:120:THR:HG23	1.60	0.82
1:B:195:THR:HG22	1:C:120:THR:CG2	2.09	0.81
1:C:110:SER:CB	1:C:129:THR:HG22	2.00	0.79
1:A:64:ILE:HD13	1:A:74:THR:HG22	1.65	0.78
1:C:110:SER:HB3	1:C:129:THR:CG2	2.00	0.77
1:A:156:LEU:O	1:A:170:GLY:HA3	1.86	0.75
1:D:94:SER:HB2	1:D:171:ARG:HG2	1.69	0.74
1:C:60:ARG:NH2	1:C:78:ASP:OD2	2.21	0.73
1:A:89:VAL:HG22	1:A:180:VAL:HG22	1.71	0.73
1:A:169:VAL:HG11	1:A:225:GLY:HA2	1.71	0.73
1:D:117:THR:HG22	1:D:185:ALA:CB	2.18	0.72
1:D:117:THR:CG2	1:D:185:ALA:CA	2.51	0.70
1:C:224:SER:HB3	1:C:230:GLY:HA2	1.71	0.70
1:D:25:ILE:HD12	1:D:65:VAL:HG12	1.73	0.69
1:D:117:THR:CB	1:D:185:ALA:HA	2.23	0.69
1:C:3:THR:C	1:C:4:ILE:HG12	2.13	0.69
1:A:64:ILE:CD1	1:A:74:THR:HG22	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:ILE:HD11	1:D:72:SER:OG	1.94	0.67
1:B:158:ARG:HH11	1:B:158:ARG:CG	2.06	0.66
1:A:193:THR:HG21	1:D:120:THR:OG1	1.96	0.66
1:B:127:HIS:CE1	1:B:129:THR:HG22	2.31	0.65
1:C:114:LYS:HB2	1:C:114:LYS:NZ	2.13	0.64
1:B:64:ILE:HG12	1:B:74:THR:HG22	1.80	0.64
1:B:195:THR:CG2	1:C:120:THR:HG23	2.22	0.64
1:C:62:SER:HB3	1:C:76:SER:HB3	1.79	0.63
1:B:3:THR:CG2	1:B:215:ASN:HD21	2.11	0.63
1:C:23:GLN:OE1	1:C:42:VAL:HG23	2.00	0.61
1:C:4:ILE:HD13	1:C:214:ALA:HB2	1.83	0.60
1:B:156:LEU:O	1:B:170:GLY:HA3	2.01	0.60
1:B:158:ARG:HG3	1:B:158:ARG:NH1	2.01	0.60
1:B:62:SER:HB3	1:B:76:SER:HB3	1.85	0.58
1:B:56:SER:HG	1:B:188:ALA:H	1.48	0.58
1:C:91:VAL:HG12	1:C:213:ILE:HG12	1.85	0.57
1:A:198:ILE:HD11	1:A:209:ILE:HD11	1.86	0.57
1:D:115:LEU:HD22	1:D:117:THR:HG23	1.85	0.57
1:D:66:SER:HB3	1:D:72:SER:HB2	1.87	0.57
1:A:171:ARG:HG3	1:A:220:ILE:HG12	1.87	0.56
1:B:49:THR:HG23	1:C:120:THR:HG22	1.88	0.56
1:B:116:LYS:HB2	1:B:187:VAL:HB	1.87	0.56
1:B:62:SER:CB	1:B:76:SER:HB3	2.36	0.56
1:A:23:GLN:HE22	1:A:205:ILE:HD11	1.69	0.56
1:C:156:LEU:O	1:C:170:GLY:HA3	2.06	0.56
1:C:94:SER:HB2	1:C:171:ARG:HG2	1.88	0.55
1:A:80:ASP:OD1	1:A:82:ASN:HB2	2.06	0.55
1:B:23:GLN:OE1	1:B:42:VAL:HG23	2.07	0.55
1:A:193:THR:HG21	1:D:120:THR:CB	2.37	0.55
1:A:94:SER:HB2	1:A:171:ARG:HG2	1.89	0.55
1:C:123:ALA:HB1	1:D:132:GLN:HG2	1.87	0.55
1:B:110:SER:OG	1:B:129:THR:HB	2.08	0.54
1:C:9:LEU:HD11	1:C:65:VAL:HG11	1.89	0.54
1:B:2:ASP:HB3	1:B:4:ILE:CD1	2.38	0.54
1:A:187:VAL:CG2	1:D:64:ILE:HG21	2.39	0.53
1:C:64:ILE:HD12	1:C:74:THR:HG22	1.91	0.52
1:A:60:ARG:HD3	1:A:76:SER:OG	2.10	0.52
1:D:7:VAL:HG21	1:D:52:ILE:HG12	1.92	0.52
1:C:117:THR:CG2	1:C:185:ALA:HA	2.39	0.52
1:C:14:ASN:HB2	1:C:17:ILE:HG12	1.91	0.51
1:D:144:GLY:C	1:D:146:ALA:H	2.12	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:GLN:HG3	1:C:125:SER:N	2.25	0.51
1:D:49:THR:HG23	1:D:195:THR:HG22	1.93	0.51
1:D:69:GLY:HA2	1:D:70:GLY:C	2.31	0.51
1:C:127:HIS:NE2	1:C:129:THR:HG23	2.26	0.51
1:B:183:LYS:C	1:B:185:ALA:H	2.13	0.51
1:D:89:VAL:HG22	1:D:180:VAL:HG23	1.92	0.50
1:A:87:GLU:HG3	1:A:181:TRP:O	2.12	0.50
1:C:117:THR:HG22	1:C:185:ALA:HA	1.93	0.50
1:B:87:GLU:HG2	1:B:88:TRP:CD1	2.46	0.49
1:A:115:LEU:HD11	1:A:181:TRP:HA	1.93	0.49
1:B:28:ASN:HB3	1:B:31:SER:O	2.12	0.49
1:D:160:SER:HB2	1:D:165:GLN:NE2	2.28	0.49
1:D:86:PRO:HG2	1:D:89:VAL:HG12	1.94	0.48
1:A:64:ILE:HG21	1:D:187:VAL:HG22	1.95	0.48
1:B:22:TYR:CD1	1:B:39:ARG:HG3	2.48	0.48
1:D:40:TRP:CZ3	1:D:42:VAL:HG22	2.49	0.48
1:D:57:VAL:HG23	1:D:187:VAL:HG13	1.95	0.48
1:D:156:LEU:O	1:D:170:GLY:HA3	2.14	0.47
1:D:22:TYR:HD1	1:D:23:GLN:O	1.96	0.47
1:A:16:ASP:OD2	1:A:227:ARG:NH1	2.46	0.47
1:D:34:SER:HB2	1:D:37:THR:HG23	1.97	0.47
1:A:32:ILE:HD11	1:A:229:LEU:HD12	1.96	0.47
1:C:23:GLN:OE1	1:C:42:VAL:CG2	2.62	0.47
1:A:198:ILE:HD11	1:A:209:ILE:CD1	2.44	0.47
1:C:138:LYS:O	1:D:177:PRO:HG2	2.14	0.47
1:C:224:SER:O	1:C:228:LEU:HB2	2.15	0.47
1:B:2:ASP:HB3	1:B:4:ILE:HD13	1.96	0.47
1:B:111:PHE:HB3	1:B:128:PHE:CE2	2.50	0.46
1:C:143:GLN:HB3	1:C:220:ILE:HD11	1.96	0.46
1:C:80:ASP:OD1	1:C:82:ASN:HB2	2.16	0.46
1:C:88:TRP:CE3	1:D:138:LYS:HB2	2.51	0.46
1:B:137:PRO:HB2	1:B:140:LEU:HB2	1.97	0.46
1:C:118:ASN:C	1:C:120:THR:H	2.19	0.46
1:B:60:ARG:HG3	1:B:78:ASP:OD1	2.14	0.46
1:A:138:LYS:O	1:B:177:PRO:HG2	2.17	0.46
1:C:89:VAL:HG22	1:C:180:VAL:HG23	1.97	0.46
1:C:77:TYR:CG	1:C:78:ASP:N	2.83	0.45
1:A:94:SER:HB3	1:A:171:ARG:HD2	1.97	0.45
1:C:131:ASN:HB2	1:D:122:ASP:O	2.16	0.45
1:A:89:VAL:HG22	1:A:180:VAL:CG2	2.43	0.45
1:C:118:ASN:C	1:C:120:THR:N	2.67	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:THR:HG21	1:D:120:THR:HB	1.99	0.44
1:C:66:SER:HB3	1:C:72:SER:HB2	1.99	0.44
1:A:11:THR:HG23	1:A:208:GLY:HA2	1.99	0.44
1:A:117:THR:HG23	1:A:118:ASN:N	2.31	0.44
1:B:70:GLY:O	1:B:71:SER:C	2.56	0.44
1:B:224:SER:HB3	1:B:232:PHE:O	2.18	0.44
1:A:34:SER:HB2	1:A:37:THR:HG23	1.99	0.44
1:A:60:ARG:CD	1:A:76:SER:OG	2.66	0.44
1:D:91:VAL:HG11	1:D:111:PHE:CZ	2.53	0.44
1:C:225:GLY:O	1:C:226:GLY:C	2.56	0.44
1:C:225:GLY:O	1:C:227:ARG:N	2.51	0.44
1:B:12:TYR:HA	1:B:13:PRO:HD3	1.81	0.43
1:D:34:SER:HB2	1:D:37:THR:CG2	2.48	0.43
1:B:93:LEU:HD13	1:B:106:ILE:HG12	2.00	0.43
1:A:206:ALA:HA	1:A:207:ASP:HA	1.91	0.43
1:A:175:TYR:CE2	1:B:175:TYR:CE2	3.06	0.43
1:C:88:TRP:CD2	1:D:138:LYS:HB2	2.54	0.42
1:B:27:ILE:HD12	1:B:61:LEU:HG	2.01	0.42
1:A:64:ILE:HG21	1:D:187:VAL:CG2	2.49	0.42
1:A:49:THR:HG21	1:D:119:SER:HA	2.01	0.42
1:A:91:VAL:HG12	1:A:213:ILE:HG12	2.01	0.42
1:C:12:TYR:HA	1:C:13:PRO:HD3	1.81	0.42
1:D:69:GLY:HA2	1:D:70:GLY:O	2.20	0.42
1:A:47:VAL:O	1:A:68:PRO:HD3	2.19	0.42
1:D:117:THR:HA	1:D:186:VAL:H	1.83	0.42
1:B:139:ASP:O	1:B:175:TYR:HB2	2.20	0.41
1:D:224:SER:HB2	1:D:230:GLY:HA2	2.02	0.41
1:B:163:SER:HA	1:B:164:PRO:HD2	1.72	0.41
1:A:33:ARG:HD2	1:A:236:ASN:O	2.20	0.41
1:D:25:ILE:HD13	1:D:75:VAL:HG12	2.01	0.41
1:C:163:SER:HA	1:C:164:PRO:HD3	1.96	0.41
1:B:99:LEU:HG	1:B:99:LEU:O	2.21	0.41
1:D:32:ILE:CD1	1:D:227:ARG:O	2.68	0.41
1:C:220:ILE:H	1:C:220:ILE:HG13	1.64	0.41
1:C:54:TYR:HB3	1:C:190:PHE:CE2	2.56	0.41
1:A:111:PHE:HB3	1:A:128:PHE:CZ	2.56	0.41
1:B:145:ASP:HB3	1:B:158:ARG:HG2	2.03	0.41
1:A:111:PHE:HB3	1:A:128:PHE:CE2	2.55	0.41
1:A:122:ASP:O	1:B:131:ASN:HB2	2.20	0.41
1:A:160:SER:N	1:A:163:SER:O	2.49	0.41
1:D:116:LYS:C	1:D:118:ASN:H	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:TYR:C	1:C:118:ASN:HD21	2.24	0.41
1:C:119:SER:O	1:C:120:THR:OG1	2.36	0.40
1:D:12:TYR:HA	1:D:13:PRO:HD3	1.92	0.40
1:B:71:SER:HB3	4:B:262:HOH:O	2.19	0.40
1:A:169:VAL:CG1	1:A:225:GLY:HA2	2.47	0.40
1:B:104:ASN:HD22	1:B:209:ILE:CG1	2.34	0.40
1:B:117:THR:HG23	1:B:118:ASN:N	2.36	0.40
1:D:90:ARG:HB3	1:D:173:LEU:HD22	2.04	0.40
1:B:131:ASN:O	1:B:132:GLN:OE1	2.40	0.40

All (12) 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:222:HIS:CG	1:C:1:ALA:N[1_655]	1.05	1.15
1:B:222:HIS:CE1	1:C:1:ALA:CB[1_655]	1.07	1.13
1:B:222:HIS:NE2	1:C:1:ALA:CA[1_655]	1.36	0.84
1:B:222:HIS:CD2	1:C:1:ALA:CA[1_655]	1.46	0.74
1:B:222:HIS:ND1	1:C:1:ALA:N[1_655]	1.58	0.62
1:B:222:HIS:CE1	1:C:1:ALA:CA[1_655]	1.62	0.58
1:B:222:HIS:CG	1:C:1:ALA:CA[1_655]	1.73	0.47
1:B:222:HIS:ND1	1:C:1:ALA:CA[1_655]	1.79	0.41
1:B:222:HIS:CB	1:C:1:ALA:N[1_655]	1.84	0.36
1:B:222:HIS:ND1	1:C:1:ALA:CB[1_655]	1.85	0.35
1:B:222:HIS:NE2	1:C:1:ALA:CB[1_655]	1.85	0.35
1:B:222:HIS:CD2	1:C:1:ALA:N[1_655]	1.92	0.28

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	234/236 (99%)	218 (93%)	15 (6%)	1 (0%)	39 74

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	234/236 (99%)	212 (91%)	16 (7%)	6 (3%)	7	26
1	C	234/236 (99%)	214 (92%)	17 (7%)	3 (1%)	15	46
1	D	234/236 (99%)	213 (91%)	17 (7%)	4 (2%)	11	38
All	All	936/944 (99%)	857 (92%)	65 (7%)	14 (2%)	13	42

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	221	PRO
1	D	121	ALA
1	B	184	SER
1	C	226	GLY
1	D	69	GLY
1	D	120	THR
1	B	161	ASN
1	C	225	GLY
1	D	221	PRO
1	B	71	SER
1	A	121	ALA
1	B	119	SER
1	B	167	ASN
1	B	70	GLY

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	199/199 (100%)	176 (88%)	23 (12%)	7	20
1	B	199/199 (100%)	178 (89%)	21 (11%)	8	25
1	C	199/199 (100%)	181 (91%)	18 (9%)	12	34
1	D	199/199 (100%)	176 (88%)	23 (12%)	7	20
All	All	796/796 (100%)	711 (89%)	85 (11%)	8	24

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	16	ASP
1	A	25	ILE
1	A	60	ARG
1	A	71	SER
1	A	74	THR
1	A	76	SER
1	A	82	ASN
1	A	108	SER
1	A	117	THR
1	A	120	THR
1	A	132	GLN
1	A	155	GLN
1	A	158	ARG
1	A	161	ASN
1	A	180	VAL
1	A	183	LYS
1	A	199	LYS
1	A	201	THR
1	A	204	ASP
1	A	205	ILE
1	A	207	ASP
1	A	234	ASP
1	B	3	THR
1	B	9	LEU
1	B	16	ASP
1	B	37	THR
1	B	64	ILE
1	B	66	SER
1	B	71	SER
1	B	76	SER
1	B	110	SER
1	B	115	LEU
1	B	120	THR
1	B	129	THR
1	B	135	GLN
1	B	147	SER
1	B	158	ARG
1	B	161	ASN
1	B	169	VAL
1	B	180	VAL
1	B	183	LYS

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Mol	Chain	Res	Type
1	B	202	ASP
1	B	218	SER
1	C	14	ASN
1	C	25	ILE
1	C	66	SER
1	C	72	SER
1	C	94	SER
1	C	110	SER
1	C	114	LYS
1	C	115	LEU
1	C	117	THR
1	C	118	ASN
1	C	124	GLN
1	C	135	GLN
1	C	175	TYR
1	C	181	TRP
1	C	183	LYS
1	C	199	LYS
1	C	207	ASP
1	C	234	ASP
1	D	25	ILE
1	D	29	ILE
1	D	32	ILE
1	D	64	ILE
1	D	72	SER
1	D	74	THR
1	D	90	ARG
1	D	101	LYS
1	D	115	LEU
1	D	116	LYS
1	D	119	SER
1	D	122	ASP
1	D	125	SER
1	D	129	THR
1	D	136	ASN
1	D	166	SER
1	D	183	LYS
1	D	201	THR
1	D	204	ASP
1	D	205	ILE
1	D	207	ASP
1	D	215	ASN

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Mol	Chain	Res	Type
1	D	222	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	118	ASN
1	A	155	GLN
1	B	82	ASN
1	B	104	ASN
1	B	131	ASN
1	B	135	GLN
1	B	161	ASN
1	B	215	ASN
1	C	118	ASN
1	C	124	GLN
1	C	155	GLN
1	D	165	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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 ⓘ

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 ⓘ

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	236/236 (100%)	-0.33	1 (0%) 93 92	19, 32, 44, 48	0
1	B	236/236 (100%)	-0.18	6 (2%) 61 55	20, 35, 50, 54	0
1	C	236/236 (100%)	-0.12	3 (1%) 79 78	23, 38, 52, 57	0
1	D	236/236 (100%)	-0.25	2 (0%) 87 86	20, 34, 44, 51	0
All	All	944/944 (100%)	-0.22	12 (1%) 79 78	19, 35, 48, 57	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	120	THR	4.4
1	B	99	LEU	3.9
1	C	119	SER	2.9
1	B	222	HIS	2.9
1	B	203	SER	2.7
1	C	118	ASN	2.6
1	B	201	THR	2.2
1	B	202	ASP	2.2
1	D	119	SER	2.1
1	D	168	SER	2.0
1	A	203	SER	2.0
1	B	162	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CA	A	237	1/1	0.97	0.07	-2.33	30,30,30,30	0
2	CA	B	241	1/1	0.96	0.06	-2.45	37,37,37,37	0
3	MN	C	244	1/1	0.98	0.10	-2.89	34,34,34,34	0
2	CA	D	239	1/1	0.97	0.07	-3.13	32,32,32,32	0
3	MN	B	242	1/1	0.97	0.08	-3.16	32,32,32,32	0
3	MN	D	240	1/1	0.97	0.07	-3.84	45,45,45,45	0
3	MN	A	238	1/1	0.99	0.08	-4.16	21,21,21,21	0
2	CA	C	243	1/1	0.93	0.06	-4.21	49,49,49,49	0

6.5 Other polymers

There are no such residues in this entry.