



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:42 AM UTC

PDB ID : 1LU2 / pdb_00001lu2
Title : DOLICHOS BIFLORUS SEED LECTIN IN COMPLEX WITH THE BLOOD GROUP A TRISACCHARIDE
Authors : Hamelryck, T.W.; Loris, R.; Bouckaert, J.; Strecker, G.; Imberty, A.; Fernandez, E.; Wyns, L.; Etzler, M.E.
Deposited on : 1998-07-30
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

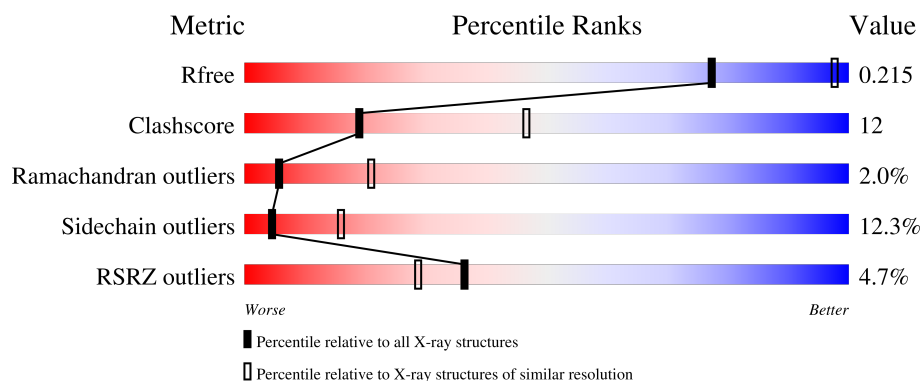
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	253	5% 72% 21% 7%
1	B	253	5% 71% 21% 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3788 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.

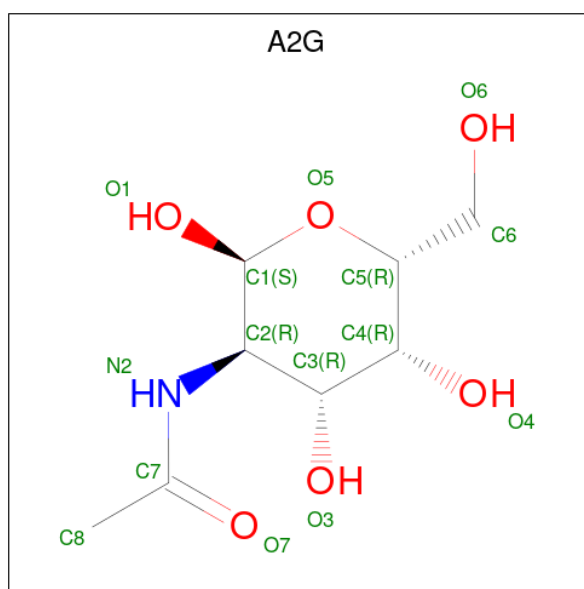
- Molecule 1 is a protein called LECTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	0	0	0
			1877	1193	302	381	1			
1	B	253	Total	C	N	O	S	0	0	0
			1877	1193	302	381	1			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	LEU	PHE	conflict	UNP P05045
B	127	LEU	PHE	conflict	UNP P05045

- Molecule 2 is 2-acetamido-2-deoxy- α -D-galactopyranose (CCD ID: A2G) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	8	1	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			15	8	1	6		

- 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	B	1	Total	Ca	0	0
			1	1		

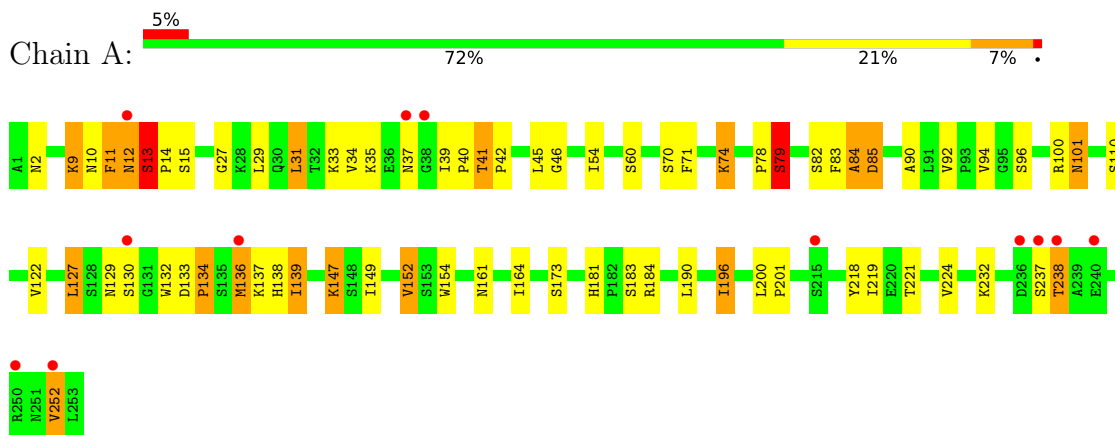
- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		

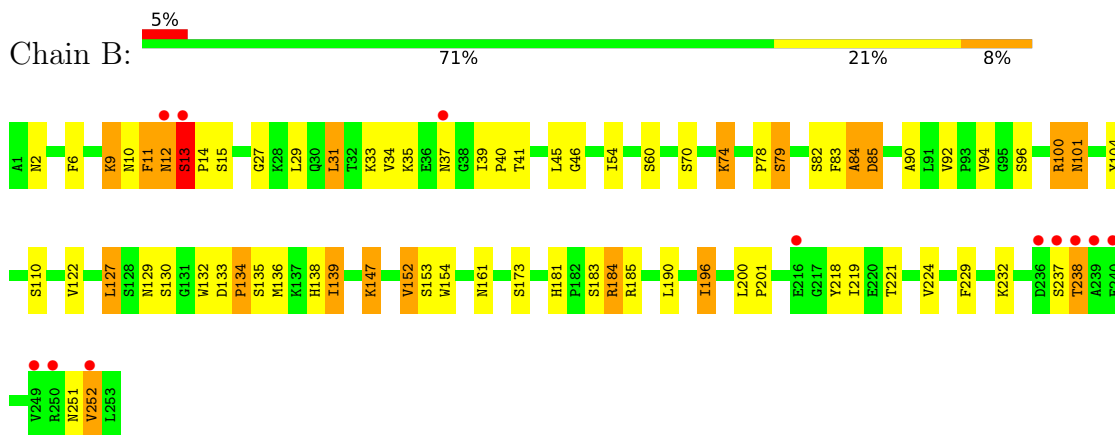
3 Residue-property plots

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: LECTIN



• Molecule 1: LECTIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	96.73Å 108.00Å 80.96Å 90.00° 124.67° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 15.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	85.3 (15.00-2.80) 84.6 (15.00-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 2.81Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.191 , 0.219 0.188 , 0.215	Depositor DCC
R_{free} test set	1457 reflections (10.17%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	1.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 60.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.208 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3788	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.60% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MN, A2G

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	0.54	0/1920	1.08	15/2628 (0.6%)
1	B	0.55	0/1920	1.07	15/2628 (0.6%)
All	All	0.54	0/3840	1.08	30/5256 (0.6%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	11	PHE	N-CA-C	13.61	130.12	113.38
1	A	11	PHE	N-CA-C	13.38	129.96	113.50
1	A	94	VAL	N-CA-C	9.10	119.91	110.36
1	A	13	SER	CA-C-N	8.73	130.14	120.45
1	A	13	SER	C-N-CA	8.73	130.14	120.45
1	B	13	SER	CA-C-N	8.37	129.75	120.45
1	B	13	SER	C-N-CA	8.37	129.75	120.45
1	A	137	LYS	N-CA-C	-6.94	97.94	109.46
1	B	252	VAL	N-CA-C	6.93	117.04	110.74
1	A	252	VAL	N-CA-C	6.91	117.03	110.74
1	B	13	SER	N-CA-C	6.72	124.66	109.81
1	A	139	ILE	N-CA-C	-6.70	98.83	108.48
1	A	13	SER	N-CA-C	6.68	124.57	109.81
1	B	251	ASN	N-CA-C	6.60	121.17	112.72
1	A	9	LYS	N-CA-C	-6.60	105.53	113.97
1	B	139	ILE	N-CA-C	-6.46	99.18	108.48
1	B	9	LYS	N-CA-C	-6.32	105.88	113.97
1	B	135	SER	N-CA-C	6.14	118.77	111.71
1	A	122	VAL	N-CA-C	-5.65	99.60	107.80
1	A	134	PRO	N-CA-C	-5.42	103.06	111.13
1	B	134	PRO	N-CA-C	-5.41	103.07	111.13
1	B	122	VAL	N-CA-C	-5.32	100.08	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	VAL	N-CA-C	5.19	113.60	107.84
1	A	92	VAL	N-CA-C	5.19	113.60	107.84
1	B	31	LEU	N-CA-C	5.12	117.76	111.82
1	A	90	ALA	N-CA-C	5.06	117.37	109.52
1	B	94	VAL	N-CA-C	5.04	119.82	109.34
1	B	90	ALA	N-CA-C	5.03	117.32	109.52
1	A	149	ILE	N-CA-C	-5.03	107.85	112.83
1	A	31	LEU	N-CA-C	5.01	117.64	111.82

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	1877	0	1792	47	0
1	B	1877	0	1792	46	0
2	A	15	0	12	1	0
2	B	15	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	3788	0	3608	87	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 12.

All (87) 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:12:ASN:O	1:A:13:SER:HB3	1.76	0.86
1:A:10:ASN:HD21	1:A:12:ASN:ND2	1.74	0.85
1:B:12:ASN:O	1:B:13:SER:HB3	1.76	0.84
1:B:10:ASN:HD21	1:B:12:ASN:ND2	1.75	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:SER:HB2	1:A:14:PRO:HD2	1.64	0.80
1:B:13:SER:HB2	1:B:14:PRO:HD2	1.64	0.78
1:A:9:LYS:H	1:B:2:ASN:ND2	1.86	0.74
1:B:11:PHE:HB2	1:B:27:GLY:O	1.90	0.72
1:A:2:ASN:ND2	1:B:9:LYS:H	1.88	0.72
1:A:11:PHE:HB2	1:A:27:GLY:O	1.90	0.71
1:A:10:ASN:HD21	1:A:12:ASN:HD21	1.40	0.69
1:A:10:ASN:ND2	1:A:12:ASN:ND2	2.42	0.68
1:A:9:LYS:H	1:B:2:ASN:HD21	1.39	0.68
1:A:12:ASN:O	1:A:13:SER:CB	2.43	0.67
1:B:10:ASN:ND2	1:B:12:ASN:ND2	2.44	0.66
1:B:12:ASN:O	1:B:13:SER:CB	2.44	0.66
1:B:10:ASN:HD21	1:B:12:ASN:HD21	1.42	0.65
1:A:2:ASN:HD21	1:B:9:LYS:H	1.43	0.63
1:A:74:LYS:NZ	1:A:161:ASN:HD21	1.96	0.62
1:B:74:LYS:NZ	1:B:161:ASN:HD21	1.98	0.61
1:B:196:ILE:HG13	1:B:200:LEU:HD12	1.83	0.60
1:B:34:VAL:HG12	1:B:35:LYS:O	2.01	0.60
1:A:134:PRO:HD2	1:A:138:HIS:CE1	2.37	0.60
1:A:196:ILE:HG13	1:A:200:LEU:HD12	1.83	0.59
1:B:136:MET:HE3	1:B:153:SER:OG	2.03	0.59
1:B:134:PRO:HD2	1:B:138:HIS:CE1	2.38	0.58
1:A:34:VAL:HG12	1:A:35:LYS:O	2.05	0.57
1:A:74:LYS:HZ3	1:A:161:ASN:HD21	1.52	0.57
1:A:134:PRO:HB2	1:A:136:MET:HE3	1.88	0.54
1:A:10:ASN:ND2	1:A:12:ASN:HD21	2.05	0.54
1:A:134:PRO:HB2	1:A:136:MET:CE	2.38	0.54
1:A:10:ASN:HD21	1:A:12:ASN:CG	2.17	0.53
1:A:13:SER:HB2	1:A:14:PRO:CD	2.37	0.52
1:B:10:ASN:HD21	1:B:12:ASN:CG	2.18	0.51
1:B:45:LEU:C	1:B:45:LEU:HD23	2.36	0.50
1:B:31:LEU:O	1:B:46:GLY:HA3	2.11	0.50
1:B:83:PHE:O	1:B:84:ALA:HB2	2.12	0.50
1:A:83:PHE:O	1:A:84:ALA:HB2	2.12	0.49
1:A:84:ALA:HB3	1:A:218:TYR:HB3	1.95	0.49
1:A:31:LEU:O	1:A:46:GLY:HA3	2.13	0.48
1:A:101:ASN:N	1:A:101:ASN:HD22	2.11	0.48
1:B:78:PRO:O	1:B:79:SER:C	2.56	0.48
1:B:13:SER:HB2	1:B:14:PRO:CD	2.37	0.48
1:B:10:ASN:ND2	1:B:12:ASN:HD21	2.08	0.48
1:B:101:ASN:HD22	1:B:101:ASN:N	2.10	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LEU:HD23	1:A:45:LEU:C	2.39	0.47
1:B:184:ARG:O	1:B:185:ARG:HB2	2.14	0.47
1:A:154:TRP:HE1	1:A:181:HIS:HE1	1.62	0.47
1:B:154:TRP:HE1	1:B:181:HIS:HE1	1.63	0.47
1:B:132:TRP:HB2	1:B:147:LYS:HG3	1.97	0.46
1:B:74:LYS:HZ3	1:B:161:ASN:HD21	1.61	0.46
1:A:70:SER:OG	1:A:252:VAL:HG22	2.16	0.45
1:B:100:ARG:HD3	1:B:104:TYR:CD2	2.52	0.45
1:A:13:SER:CB	1:A:14:PRO:HD2	2.38	0.45
1:A:2:ASN:HD21	1:B:10:ASN:H	1.65	0.45
1:A:10:ASN:H	1:B:2:ASN:HD21	1.64	0.45
1:B:84:ALA:HB3	1:B:218:TYR:HB3	1.98	0.45
1:A:74:LYS:NZ	1:A:161:ASN:ND2	2.64	0.44
1:A:13:SER:CB	1:A:14:PRO:CD	2.94	0.44
1:A:40:PRO:HB2	1:A:219:ILE:HB	2.00	0.44
1:B:70:SER:OG	1:B:252:VAL:HG22	2.17	0.44
1:A:37:ASN:ND2	1:A:39:ILE:HG23	2.33	0.44
1:A:127:LEU:HD22	2:A:254:A2G:H4	1.99	0.43
1:B:84:ALA:HA	1:B:85:ASP:HA	1.76	0.43
1:A:84:ALA:HA	1:A:85:ASP:HA	1.76	0.43
1:A:40:PRO:HG3	1:A:221:THR:CG2	2.49	0.43
1:B:29:LEU:HB3	1:B:224:VAL:HB	2.00	0.43
1:A:200:LEU:HB3	1:A:201:PRO:HD2	2.00	0.43
1:B:40:PRO:HB2	1:B:219:ILE:HB	2.00	0.43
1:A:29:LEU:HB3	1:A:224:VAL:HB	2.00	0.42
1:B:127:LEU:HD22	2:B:254:A2G:H4	2.01	0.42
1:B:200:LEU:HB3	1:B:201:PRO:HD2	2.00	0.42
1:A:11:PHE:HB2	1:A:27:GLY:C	2.44	0.42
1:B:11:PHE:HB2	1:B:27:GLY:C	2.44	0.42
1:B:40:PRO:HG3	1:B:221:THR:CG2	2.50	0.42
1:B:37:ASN:ND2	1:B:39:ILE:HG23	2.35	0.42
1:A:129:ASN:N	1:A:133:ASP:HB2	2.35	0.42
1:A:139:ILE:HB	1:A:152:VAL:HG12	2.01	0.42
1:B:129:ASN:N	1:B:133:ASP:HB2	2.35	0.42
1:B:13:SER:CB	1:B:14:PRO:HD2	2.38	0.41
1:A:78:PRO:O	1:A:79:SER:C	2.63	0.41
1:B:139:ILE:HB	1:B:152:VAL:HG12	2.01	0.41
1:B:6:PHE:CZ	1:B:229:PHE:HB3	2.56	0.41
1:B:74:LYS:HZ2	1:B:161:ASN:ND2	2.19	0.40
1:A:132:TRP:HB2	1:A:147:LYS:HG3	2.03	0.40
1:A:41:THR:HA	1:A:42:PRO:HD3	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:PHE:CE2	1:A:164:ILE:HD13	2.57	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	251/253 (99%)	232 (92%)	14 (6%)	5 (2%)	6	21
1	B	251/253 (99%)	233 (93%)	13 (5%)	5 (2%)	6	21
All	All	502/506 (99%)	465 (93%)	27 (5%)	10 (2%)	6	21

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	79	SER
1	B	13	SER
1	B	79	SER
1	A	84	ALA
1	A	96	SER
1	A	238	THR
1	B	84	ALA
1	B	96	SER
1	B	238	THR

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	204/214 (95%)	178 (87%)	26 (13%)	4	15
1	B	204/214 (95%)	180 (88%)	24 (12%)	5	17
All	All	408/428 (95%)	358 (88%)	50 (12%)	4	16

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	15	SER
1	A	33	LYS
1	A	41	THR
1	A	54	ILE
1	A	60	SER
1	A	74	LYS
1	A	79	SER
1	A	82	SER
1	A	85	ASP
1	A	100	ARG
1	A	101	ASN
1	A	110	SER
1	A	127	LEU
1	A	130	SER
1	A	136	MET
1	A	147	LYS
1	A	152	VAL
1	A	173	SER
1	A	183	SER
1	A	184	ARG
1	A	190	LEU
1	A	196	ILE
1	A	232	LYS
1	A	237	SER
1	A	238	THR
1	B	12	ASN
1	B	15	SER
1	B	33	LYS
1	B	41	THR
1	B	54	ILE
1	B	60	SER
1	B	74	LYS

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Mol	Chain	Res	Type
1	B	82	SER
1	B	85	ASP
1	B	100	ARG
1	B	101	ASN
1	B	110	SER
1	B	127	LEU
1	B	130	SER
1	B	147	LYS
1	B	152	VAL
1	B	173	SER
1	B	183	SER
1	B	184	ARG
1	B	190	LEU
1	B	196	ILE
1	B	232	LYS
1	B	237	SER
1	B	238	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	10	ASN
1	A	12	ASN
1	A	55	GLN
1	A	101	ASN
1	A	115	ASN
1	A	118	GLN
1	A	161	ASN
1	A	181	HIS
1	A	198	ASN
1	A	222	HIS
1	B	2	ASN
1	B	10	ASN
1	B	12	ASN
1	B	30	GLN
1	B	55	GLN
1	B	101	ASN
1	B	115	ASN
1	B	118	GLN
1	B	161	ASN
1	B	181	HIS

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Mol	Chain	Res	Type
1	B	198	ASN
1	B	222	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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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	A2G	B	254	-	15,15,15	0.67	0	21,21,21	0.64	0
2	A2G	A	254	-	15,15,15	0.65	0	21,21,21	0.63	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	A2G	B	254	-	-	0/6/26/26	0/1/1/1
2	A2G	A	254	-	-	0/6/26/26	0/1/1/1

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	254	A2G	1	0
2	A	254	A2G	1	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	253/253 (100%)	-0.06	12 (4%)	36	29	13, 29, 47, 66	18 (7%)
1	B	253/253 (100%)	-0.00	12 (4%)	36	29	14, 30, 46, 65	18 (7%)
All	All	506/506 (100%)	-0.03	24 (4%)	36	29	13, 29, 47, 66	36 (7%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	ASP	5.1
1	B	237	SER	4.3
1	A	240	GLU	4.0
1	B	236	ASP	4.0
1	B	240	GLU	3.7
1	A	237	SER	3.6
1	B	238	THR	3.5
1	B	12	ASN	3.2
1	B	239	ALA	3.0
1	A	12	ASN	3.0
1	B	252	VAL	2.8
1	A	238	THR	2.8
1	A	130	SER	2.7
1	A	250	ARG	2.6
1	A	215	SER	2.5
1	B	216	GLU	2.5
1	B	13	SER	2.4
1	B	249	VAL	2.4
1	A	252	VAL	2.4
1	B	250	ARG	2.3
1	B	37	ASN	2.2
1	A	38	GLY	2.1
1	A	37	ASN	2.1
1	A	136	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	A2G	B	254	15/15	0.88	0.11	54,55,58,60	0
4	MN	A	302	1/1	0.88	0.15	70,70,70,70	0
2	A2G	A	254	15/15	0.91	0.12	55,57,60,62	0
3	CA	B	301	1/1	0.95	0.04	28,28,28,28	0
4	MN	B	302	1/1	0.95	0.08	74,74,74,74	0
3	CA	A	301	1/1	0.97	0.05	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.