



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:02 AM UTC

PDB ID : 4DBA / pdb\_00004dba  
Title : Designed Armadillo repeat protein (YIIM3AII)  
Authors : Madhurantakam, C.; Varadamsetty, G.; Grutter, M.G.; Pluckthun, A.; Mittl, P.R.E.  
Deposited on : 2012-01-13  
Resolution : 2.40 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](mailto: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>.

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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

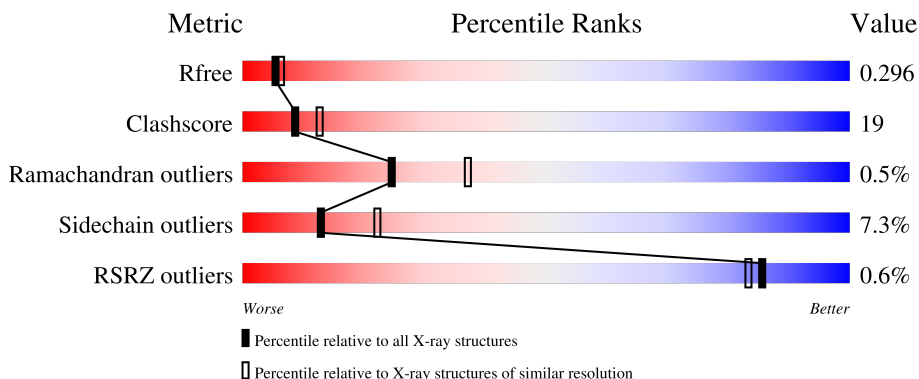
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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  $\geq 3$ , 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	210	 61% 30% 7%
1	B	210	 63% 26% 7%
1	C	210	 60% 28% 5% 7%
1	D	210	 60% 30% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	B	302	-	-	X	-

## 2 Entry composition [i](#)

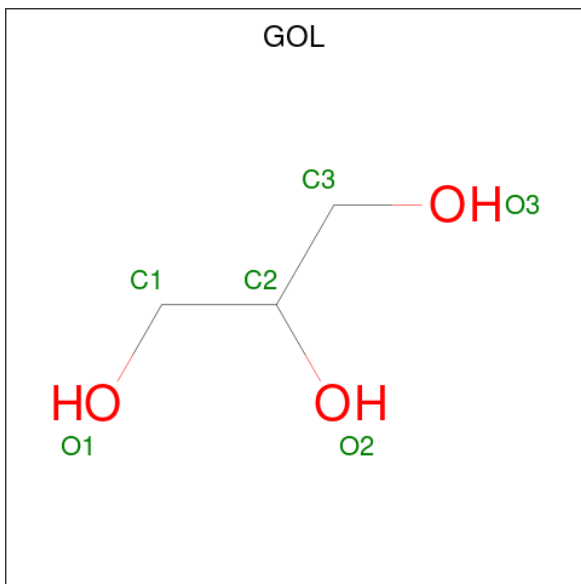
There are 3 unique types of molecules in this entry. The entry contains 5976 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Designed Armadillo repeat protein, YIIM3AII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	196	Total 1470	C 912	N 251	O 305	S 2	0	0	0
1	B	196	Total 1466	C 910	N 251	O 303	S 2	0	0	0
1	C	196	Total 1470	C 912	N 251	O 305	S 2	0	0	0
1	D	196	Total 1470	C 912	N 251	O 305	S 2	0	0	0

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	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	13	Total	O	0	0
			13	13		
3	B	22	Total	O	0	0
			22	22		
3	C	21	Total	O	0	0
			21	21		
3	D	14	Total	O	0	0
			14	14		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.15Å 60.60Å 61.86Å 74.82° 89.55° 75.53°	Depositor
Resolution (Å)	25.29 – 2.40 25.29 – 2.40	Depositor EDS
% Data completeness (in resolution range)	89.7 (25.29-2.40) 89.6 (25.29-2.40)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 2.39Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.234 , 0.301 0.232 , 0.296	Depositor DCC
$R_{free}$ test set	1434 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	57.5	Xtrriage
Anisotropy	0.256	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 53.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5976	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.63	0/1484	0.93	1/2015 (0.0%)
1	B	0.65	0/1480	0.97	1/2010 (0.0%)
1	C	0.59	0/1484	0.91	0/2015
1	D	0.56	0/1484	0.95	1/2015 (0.0%)
All	All	0.61	0/5932	0.94	3/8055 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49	VAL	N-CA-C	-6.71	106.04	111.81
1	D	139	LEU	N-CA-C	5.85	120.71	112.75
1	B	49	VAL	N-CA-C	-5.44	107.13	111.81

There are no chirality outliers.

There are no planarity outliers.

### 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	1470	0	1466	52	0
1	B	1466	0	1462	53	0
1	C	1470	0	1466	63	0
1	D	1470	0	1466	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	18	0	24	4	0
2	D	12	0	16	0	0
3	A	13	0	0	7	0
3	B	22	0	0	7	0
3	C	21	0	0	10	0
3	D	14	0	0	14	0
All	All	5976	0	5900	222	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 19.

All (222) 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:15:GLN:HB2	3:D:405:HOH:O	1.60	1.00
1:B:75:TRP:HZ3	2:B:302:GOL:HO2	1.00	0.95
1:B:16:MET:HE2	1:B:34:LYS:HD3	1.52	0.90
1:C:152:GLN:HG3	3:C:310:HOH:O	1.73	0.88
1:D:185:GLU:O	1:D:188:GLN:HG2	1.79	0.83
1:A:32:THR:CG2	1:B:72:GLU:HB2	2.10	0.80
1:C:86:GLU:OE1	1:C:86:GLU:HA	1.83	0.78
1:C:168:GLY:O	1:C:172:LYS:HG3	1.85	0.76
1:B:169:ASN:HA	1:B:172:LYS:HE3	1.68	0.75
1:C:147:SER:HB3	1:C:183:LYS:NZ	2.01	0.75
1:B:101:VAL:O	1:B:104:LEU:HB2	1.88	0.74
1:A:101:VAL:O	1:A:104:LEU:HB2	1.88	0.73
1:C:72:GLU:HB3	1:D:32:THR:HG22	1.71	0.73
1:C:72:GLU:HB3	3:D:413:HOH:O	1.88	0.73
1:C:28:GLN:NE2	1:D:61:LEU:HD21	2.04	0.72
1:C:183:LYS:HD2	1:C:186:GLN:NE2	2.06	0.71
1:B:183:LYS:HE3	3:B:415:HOH:O	1.91	0.70
1:C:12:GLU:C	1:C:14:PRO:HD2	2.17	0.69
1:B:131:GLN:NE2	1:B:171:GLN:OE1	2.26	0.69
1:D:59:VAL:HG13	3:D:411:HOH:O	1.93	0.68
1:A:89:GLN:HG2	1:A:129:GLN:HE21	1.58	0.68
1:A:32:THR:HG22	1:B:72:GLU:HB2	1.75	0.67
1:D:139:LEU:O	1:D:143:VAL:HG23	1.93	0.67
1:C:201:GLU:O	1:C:205:LYS:HG3	1.94	0.67
1:A:16:MET:HE2	1:A:34:LYS:HD3	1.75	0.67
1:A:23:ASP:HB2	3:A:306:HOH:O	1.94	0.67
1:A:32:THR:HG21	1:B:72:GLU:HB2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:GLU:O	1:B:188:GLN:HG2	1.94	0.66
1:A:89:GLN:HG2	1:A:129:GLN:NE2	2.10	0.66
1:D:67:GLU:CD	1:D:67:GLU:H	2.03	0.66
1:B:204:GLU:O	1:B:207:GLN:HG3	1.96	0.66
1:C:13:LEU:N	1:C:14:PRO:HD2	2.10	0.66
1:D:105:SER:HA	3:D:408:HOH:O	1.96	0.65
1:C:16:MET:HE2	1:C:34:LYS:HD3	1.79	0.64
1:D:134:ILE:HG12	1:D:139:LEU:HD11	1.78	0.64
1:A:95:GLY:HA2	3:A:301:HOH:O	1.97	0.64
1:C:171:GLN:O	1:C:175:VAL:HG23	1.98	0.64
1:B:87:GLN:HG3	3:B:410:HOH:O	1.98	0.63
1:A:108:ASN:OD1	1:A:110:GLN:HB2	1.99	0.63
1:A:14:PRO:HA	3:A:313:HOH:O	1.98	0.63
1:B:78:SER:HB2	1:B:118:ALA:HB2	1.78	0.63
1:C:28:GLN:HE21	1:D:61:LEU:HD21	1.63	0.63
1:C:151:GLU:O	1:C:155:GLN:HG3	2.00	0.61
1:B:12:GLU:N	3:B:413:HOH:O	2.33	0.60
1:D:150:ASN:OD1	1:D:152:GLN:HG2	2.01	0.60
1:C:176:LYS:HA	3:C:317:HOH:O	2.01	0.60
1:D:197:LYS:O	1:D:200:GLN:HB2	2.02	0.60
1:C:147:SER:HB3	1:C:183:LYS:HZ3	1.66	0.59
1:A:140:PRO:HA	3:A:309:HOH:O	2.01	0.59
1:A:143:VAL:HB	3:A:309:HOH:O	2.01	0.59
1:C:89:GLN:HG2	1:C:129:GLN:HE21	1.67	0.59
1:C:180:ALA:O	1:C:184:LEU:HB2	2.03	0.59
2:B:301:GOL:H2	2:B:302:GOL:O3	2.03	0.58
1:A:67:GLU:CD	1:A:67:GLU:H	2.10	0.58
1:C:104:LEU:HB2	3:C:316:HOH:O	2.03	0.58
1:C:72:GLU:HB3	1:D:32:THR:CG2	2.32	0.58
1:B:68:GLN:O	1:B:72:GLU:HG2	2.03	0.58
1:D:101:VAL:O	1:D:104:LEU:HB2	2.03	0.58
1:B:37:GLN:HA	1:B:37:GLN:NE2	2.19	0.57
1:A:72:GLU:HG2	3:A:312:HOH:O	2.03	0.57
1:D:32:THR:HG22	3:D:413:HOH:O	2.05	0.57
1:D:99:ALA:CB	3:D:411:HOH:O	2.52	0.57
1:A:134:ILE:HG12	1:A:139:LEU:HD11	1.87	0.57
1:C:147:SER:HB3	1:C:183:LYS:CE	2.35	0.57
1:C:188:GLN:OE1	1:C:200:GLN:HG3	2.04	0.57
1:D:167:GLY:HA3	1:D:171:GLN:HG3	1.86	0.56
1:B:206:LEU:O	1:B:207:GLN:HB2	2.03	0.56
1:B:139:LEU:O	1:B:143:VAL:HG23	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:LEU:N	1:C:14:PRO:CD	2.68	0.56
1:D:186:GLN:C	1:D:188:GLN:H	2.14	0.56
1:A:72:GLU:HB3	1:B:32:THR:HG22	1.87	0.56
1:B:37:GLN:HA	1:B:37:GLN:HE21	1.70	0.56
1:B:75:TRP:HZ3	2:B:302:GOL:O2	1.78	0.56
1:C:101:VAL:HG13	3:C:316:HOH:O	2.06	0.55
1:A:128:GLU:CD	1:A:128:GLU:H	2.15	0.55
1:C:55:LEU:HD11	1:C:80:ILE:HD13	1.89	0.55
1:D:14:PRO:HD2	3:D:404:HOH:O	2.07	0.55
1:D:185:GLU:HA	1:D:203:LEU:HD11	1.89	0.55
1:C:150:ASN:HB3	1:C:153:ILE:HD12	1.88	0.55
1:A:55:LEU:HD11	1:A:80:ILE:HD13	1.89	0.54
1:D:197:LYS:HB2	1:D:197:LYS:NZ	2.22	0.54
1:A:201:GLU:O	1:A:205:LYS:HG3	2.08	0.53
1:D:67:GLU:O	1:D:71:GLN:HG3	2.09	0.53
1:D:197:LYS:HA	1:D:200:GLN:OE1	2.09	0.53
1:C:67:GLU:O	1:C:71:GLN:HG3	2.09	0.53
1:C:176:LYS:C	1:C:178:ALA:H	2.17	0.53
1:B:89:GLN:HG2	1:B:129:GLN:HE21	1.74	0.52
1:C:74:LEU:HG	1:C:114:GLU:HG3	1.90	0.52
1:C:172:LYS:HB3	1:C:206:LEU:HD12	1.90	0.52
1:D:152:GLN:O	1:D:155:GLN:HB2	2.10	0.52
1:D:207:GLN:HB2	3:D:403:HOH:O	2.09	0.52
1:A:139:LEU:O	1:A:143:VAL:HG23	2.10	0.52
1:B:22:SER:O	1:B:28:GLN:NE2	2.43	0.51
1:B:158:LEU:HG	1:B:198:GLU:HG2	1.92	0.51
1:C:75:TRP:CH2	1:C:79:ASN:ND2	2.78	0.51
1:C:202:ALA:O	1:C:206:LEU:HD22	2.10	0.51
1:D:62:LEU:HD12	3:D:411:HOH:O	2.09	0.51
1:A:112:LEU:C	1:A:112:LEU:HD23	2.36	0.51
1:A:50:ILE:HG23	1:A:55:LEU:HD22	1.93	0.51
1:D:67:GLU:CD	1:D:67:GLU:N	2.69	0.51
1:B:181:LEU:HD22	3:B:417:HOH:O	2.11	0.50
1:C:70:LEU:HD22	1:C:111:ILE:HG12	1.93	0.50
1:C:104:LEU:HD12	3:C:316:HOH:O	2.12	0.50
1:D:197:LYS:HB2	1:D:197:LYS:HZ3	1.77	0.50
1:B:195:ILE:O	1:B:196:GLN:C	2.53	0.50
1:C:113:GLN:NE2	1:C:152:GLN:HB3	2.26	0.50
1:D:95:GLY:HA2	3:D:407:HOH:O	2.11	0.50
1:A:190:HIS:O	1:A:191:GLU:C	2.55	0.49
1:C:139:LEU:HB2	3:C:303:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:GLN:O	1:A:177:GLU:HB2	2.12	0.49
1:D:101:VAL:HG22	1:D:138:ALA:HA	1.93	0.49
1:C:78:SER:HB3	1:C:118:ALA:HB2	1.95	0.49
1:B:72:GLU:HB3	3:B:406:HOH:O	2.13	0.49
1:D:163:ASN:O	1:D:166:SER:HB2	2.12	0.49
1:B:139:LEU:HB2	1:B:140:PRO:HD3	1.95	0.49
1:D:22:SER:OG	1:D:23:ASP:N	2.46	0.49
1:D:192:ASN:HB3	1:D:195:ILE:HD12	1.94	0.48
1:C:89:GLN:HG2	1:C:129:GLN:NE2	2.28	0.48
1:D:150:ASN:OD1	1:D:150:ASN:C	2.56	0.48
1:A:179:GLY:O	1:A:180:ALA:C	2.56	0.48
1:C:19:GLN:O	1:C:22:SER:HB3	2.13	0.48
1:C:32:THR:HG22	1:D:72:GLU:HB3	1.96	0.48
1:D:16:MET:HG3	3:D:410:HOH:O	2.14	0.47
1:C:70:LEU:C	1:C:70:LEU:HD23	2.40	0.47
1:C:59:VAL:HG11	1:C:95:GLY:O	2.13	0.47
1:B:168:GLY:H	1:B:171:GLN:CG	2.26	0.47
1:D:116:LEU:HD13	1:D:156:GLU:HB3	1.95	0.47
1:C:17:THR:O	1:C:21:ASN:ND2	2.48	0.47
1:A:116:LEU:HD11	1:A:145:LEU:HD13	1.97	0.47
2:B:301:GOL:H2	2:B:302:GOL:C3	2.45	0.47
1:A:192:ASN:HD21	1:A:194:LYS:HB2	1.81	0.46
1:C:112:LEU:C	1:C:112:LEU:HD23	2.40	0.46
1:C:206:LEU:O	1:C:207:GLN:CB	2.64	0.46
1:D:13:LEU:N	1:D:14:PRO:CD	2.79	0.46
1:D:156:GLU:O	1:D:157:ALA:C	2.58	0.46
1:A:134:ILE:HD11	1:A:164:ILE:CG2	2.46	0.46
1:D:17:THR:O	1:D:21:ASN:ND2	2.48	0.46
1:D:20:LEU:HD11	1:D:35:PHE:CE2	2.50	0.46
1:A:193:GLU:HG2	1:A:194:LYS:N	2.30	0.46
1:B:77:LEU:HD12	1:B:77:LEU:HA	1.76	0.46
1:B:167:GLY:HA3	1:B:171:GLN:HG3	1.96	0.46
1:C:34:LYS:HE3	3:C:302:HOH:O	2.14	0.46
1:A:22:SER:OG	1:A:23:ASP:N	2.49	0.46
1:B:198:GLU:HG3	3:B:412:HOH:O	2.15	0.46
1:A:143:VAL:O	1:A:146:LEU:HB2	2.16	0.45
1:B:16:MET:HG2	3:B:402:HOH:O	2.16	0.45
1:B:168:GLY:H	1:B:171:GLN:HG3	1.82	0.45
1:D:156:GLU:O	1:D:159:TRP:N	2.50	0.45
1:A:15:GLN:HB3	3:A:302:HOH:O	2.17	0.45
1:A:97:LEU:O	1:A:98:PRO:C	2.58	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:GLY:N	1:B:171:GLN:CG	2.80	0.45
1:B:17:THR:HG22	1:B:17:THR:O	2.16	0.45
1:C:139:LEU:O	1:C:143:VAL:HG23	2.17	0.45
1:A:19:GLN:O	1:A:22:SER:HB3	2.17	0.45
1:C:190:HIS:CE1	1:C:195:ILE:HD12	2.52	0.45
1:C:87:GLN:N	1:C:87:GLN:OE1	2.50	0.45
1:D:99:ALA:HB3	3:D:411:HOH:O	2.17	0.45
1:D:152:GLN:HG2	1:D:152:GLN:H	1.50	0.45
1:B:12:GLU:C	1:B:14:PRO:HD2	2.42	0.44
1:C:153:ILE:O	1:C:154:LEU:C	2.60	0.44
1:B:180:ALA:O	1:B:184:LEU:HB2	2.18	0.44
1:C:147:SER:HA	3:C:306:HOH:O	2.16	0.44
1:A:81:ALA:O	1:A:121:ASN:HB3	2.16	0.44
1:A:185:GLU:HA	1:A:203:LEU:HD11	1.99	0.44
1:B:181:LEU:O	1:B:185:GLU:HB2	2.18	0.44
1:D:206:LEU:O	1:D:207:GLN:HB3	2.17	0.44
1:B:154:LEU:HD23	1:B:154:LEU:O	2.17	0.44
1:A:130:ILE:CG2	1:A:164:ILE:HG23	2.48	0.44
1:A:150:ASN:HB3	1:A:153:ILE:HD12	1.99	0.44
1:D:134:ILE:HG12	1:D:139:LEU:CD1	2.47	0.44
1:D:201:GLU:O	1:D:205:LYS:HG3	2.17	0.44
1:A:12:GLU:C	1:A:14:PRO:HD2	2.43	0.44
1:A:134:ILE:HG23	1:A:139:LEU:HD12	2.00	0.44
1:A:192:ASN:ND2	1:A:194:LYS:HB2	2.33	0.44
1:C:197:LYS:O	1:C:200:GLN:HB2	2.17	0.44
1:D:186:GLN:C	1:D:188:GLN:N	2.76	0.44
1:B:112:LEU:HD23	1:B:112:LEU:C	2.43	0.43
1:C:72:GLU:CB	1:D:32:THR:CG2	2.96	0.43
1:C:207:GLN:HB2	3:C:315:HOH:O	2.17	0.43
1:D:180:ALA:O	1:D:184:LEU:HB2	2.19	0.43
1:D:184:LEU:HD23	1:D:203:LEU:HD23	2.00	0.43
1:C:72:GLU:HG2	3:C:319:HOH:O	2.18	0.43
1:A:13:LEU:N	1:A:14:PRO:CD	2.81	0.43
1:A:116:LEU:HD23	1:A:116:LEU:HA	1.87	0.43
1:A:154:LEU:HD13	1:A:195:ILE:HD13	2.00	0.43
1:A:127:ASN:O	1:A:171:GLN:NE2	2.52	0.43
1:D:133:VAL:HG11	1:D:164:ILE:HD11	2.00	0.43
1:A:134:ILE:HG12	1:A:139:LEU:CD1	2.47	0.43
1:B:12:GLU:O	1:B:16:MET:HG3	2.19	0.43
1:B:113:GLN:NE2	1:B:152:GLN:HB3	2.34	0.43
1:C:139:LEU:HB2	1:C:140:PRO:HD3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:GLY:HA3	3:D:406:HOH:O	2.19	0.42
1:D:150:ASN:OD1	1:D:152:GLN:CG	2.67	0.42
1:A:147:SER:HB3	1:A:183:LYS:CE	2.48	0.42
1:B:24:ASP:HB3	1:B:27:GLU:HG3	2.01	0.42
1:D:74:LEU:HD21	1:D:114:GLU:HB3	2.01	0.42
1:A:197:LYS:HE2	1:A:201:GLU:OE1	2.20	0.42
1:B:148:SER:OG	1:B:149:PRO:HD2	2.19	0.42
1:B:168:GLY:N	1:B:171:GLN:HG2	2.34	0.42
1:D:203:LEU:HD23	1:D:203:LEU:HA	1.79	0.42
1:B:70:LEU:HD23	1:B:70:LEU:C	2.44	0.42
1:D:108:ASN:OD1	1:D:108:ASN:C	2.62	0.42
1:C:72:GLU:CB	1:D:32:THR:HG21	2.50	0.42
1:A:108:ASN:HB3	1:A:111:ILE:HD12	2.01	0.42
1:B:104:LEU:HD23	1:B:104:LEU:HA	1.65	0.42
1:C:165:ALA:HB1	1:C:206:LEU:HD13	2.02	0.42
1:B:207:GLN:HB2	1:B:207:GLN:HE21	1.66	0.42
1:C:32:THR:HG21	1:D:72:GLU:HB2	2.02	0.41
1:D:24:ASP:HB3	1:D:27:GLU:HG2	2.02	0.41
1:A:88:ILE:O	1:A:92:ILE:HG13	2.20	0.41
1:B:146:LEU:HB2	1:B:183:LYS:HG2	2.02	0.41
1:C:176:LYS:C	1:C:178:ALA:N	2.78	0.41
1:D:16:MET:HE2	1:D:34:LYS:HD3	2.01	0.41
1:B:15:GLN:HG2	1:B:16:MET:N	2.30	0.41
1:D:26:GLN:O	1:D:30:SER:HB2	2.20	0.41
1:C:108:ASN:OD1	1:C:110:GLN:HB2	2.20	0.41
1:D:99:ALA:O	1:D:102:GLN:HB3	2.21	0.41
1:D:147:SER:HA	3:D:412:HOH:O	2.20	0.41
1:A:154:LEU:O	1:A:158:LEU:HB2	2.21	0.41
1:C:17:THR:O	1:C:17:THR:HG22	2.20	0.40
1:B:108:ASN:OD1	1:B:110:GLN:HB2	2.22	0.40
1:B:153:ILE:O	1:B:154:LEU:C	2.64	0.40

There are no symmetry-related clashes.

## 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	194/210 (92%)	179 (92%)	13 (7%)	2 (1%)	12	20
1	B	194/210 (92%)	185 (95%)	8 (4%)	1 (0%)	24	37
1	C	194/210 (92%)	183 (94%)	11 (6%)	0	100	100
1	D	194/210 (92%)	173 (89%)	20 (10%)	1 (0%)	24	37
All	All	776/840 (92%)	720 (93%)	52 (7%)	4 (0%)	24	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	GLY
1	B	189	SER
1	A	22	SER
1	D	149	PRO

### 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	157/170 (92%)	149 (95%)	8 (5%)	21	37
1	B	156/170 (92%)	143 (92%)	13 (8%)	10	17
1	C	157/170 (92%)	145 (92%)	12 (8%)	12	21
1	D	157/170 (92%)	144 (92%)	13 (8%)	10	17
All	All	627/680 (92%)	581 (93%)	46 (7%)	13	22

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	GLU
1	A	79	ASN
1	A	104	LEU
1	A	111	ILE

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	120	SER
1	A	154	LEU
1	A	181	LEU
1	A	201	GLU
1	B	15	GLN
1	B	60	GLN
1	B	72	GLU
1	B	77	LEU
1	B	120	SER
1	B	128	GLU
1	B	170	GLU
1	B	171	GLN
1	B	181	LEU
1	B	183	LYS
1	B	184	LEU
1	B	197	LYS
1	B	207	GLN
1	C	13	LEU
1	C	67	GLU
1	C	79	ASN
1	C	86	GLU
1	C	87	GLN
1	C	111	ILE
1	C	120	SER
1	C	154	LEU
1	C	170	GLU
1	C	171	GLN
1	C	197	LYS
1	C	206	LEU
1	D	40	SER
1	D	44	GLU
1	D	67	GLU
1	D	89	GLN
1	D	92	ILE
1	D	104	LEU
1	D	110	GLN
1	D	116	LEU
1	D	120	SER
1	D	152	GLN
1	D	155	GLN
1	D	171	GLN
1	D	197	LYS

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	21	ASN
1	A	68	GLN
1	A	173	GLN
1	A	186	GLN
1	B	37	GLN
1	B	43	ASN
1	B	45	GLN
1	B	71	GLN
1	B	89	GLN
1	B	131	GLN
1	B	171	GLN
1	B	207	GLN
1	C	21	ASN
1	C	28	GLN
1	C	43	ASN
1	C	79	ASN
1	C	131	GLN
1	C	186	GLN
1	C	200	GLN
1	D	28	GLN
1	D	43	ASN
1	D	79	ASN
1	D	173	GLN

### 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

5 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	B	302	-	5,5,5	0.36	0	5,5,5	0.20	0
2	GOL	B	303	-	5,5,5	0.41	0	5,5,5	0.19	0
2	GOL	D	301	-	5,5,5	0.35	0	5,5,5	0.28	0
2	GOL	D	302	-	5,5,5	0.33	0	5,5,5	0.29	0
2	GOL	B	301	-	5,5,5	0.37	0	5,5,5	0.46	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. '2' means no outliers of that kind were identified.

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

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

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

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	GOL	4	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	196/210 (93%)	0.04	3 (1%) 72 68	37, 63, 118, 129	0
1	B	196/210 (93%)	-0.05	1 (0%) 87 85	36, 64, 98, 111	0
1	C	196/210 (93%)	0.06	1 (0%) 87 85	38, 68, 113, 122	0
1	D	196/210 (93%)	-0.01	0 100 100	47, 72, 114, 124	0
All	All	784/840 (93%)	0.01	5 (0%) 85 83	36, 67, 113, 129	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	195	ILE	3.5
1	A	191	GLU	2.9
1	A	169	ASN	2.6
1	B	189	SER	2.4
1	C	195	ILE	2.3

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	GOL	B	303	6/6	0.58	0.14	87,90,93,96	0
2	GOL	D	301	6/6	0.67	0.13	84,87,91,92	0
2	GOL	B	301	6/6	0.68	0.13	79,86,92,94	0
2	GOL	B	302	6/6	0.78	0.10	76,77,81,83	0
2	GOL	D	302	6/6	0.81	0.08	89,90,94,97	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.