



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 4, 2023 – 11:30 PM EDT

PDB ID : 3TGO  
Title : Crystal structure of the E. coli BamCD complex  
Authors : Paetzel, M.; Kim, K.H.; Aulakh, S.  
Deposited on : 2011-08-17  
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](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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35

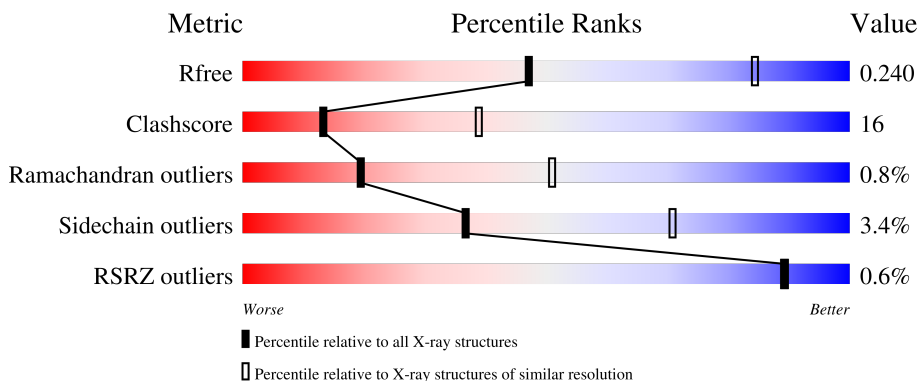
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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 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	229	 69% 25% • 5%
1	B	229	 71% 24% 5%
2	C	323	 43% 12% • 44%
2	D	323	 37% 20% • 42%

## 2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 6466 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 UPF0169 lipoprotein yfiO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	217	Total	C	N	O	S	0	0	0
			1754	1106	308	333	7			
1	B	218	Total	C	N	O	S	0	0	0
			1761	1109	309	336	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	GLY	-	expression tag	UNP P0AC02
A	18	SER	-	expression tag	UNP P0AC02
A	19	HIS	-	expression tag	UNP P0AC02
A	20	MET	-	expression tag	UNP P0AC02
B	17	GLY	-	expression tag	UNP P0AC02
B	18	SER	-	expression tag	UNP P0AC02
B	19	HIS	-	expression tag	UNP P0AC02
B	20	MET	-	expression tag	UNP P0AC02

- Molecule 2 is a protein called Lipoprotein 34.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	182	Total	C	N	O	S	0	1	0
			1383	864	238	277	4			
2	D	186	Total	C	N	O	S	0	1	0
			1412	881	244	283	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	22	GLY	-	expression tag	UNP P0A903
C	23	SER	-	expression tag	UNP P0A903
C	24	HIS	-	expression tag	UNP P0A903
C	25	MET	-	expression tag	UNP P0A903

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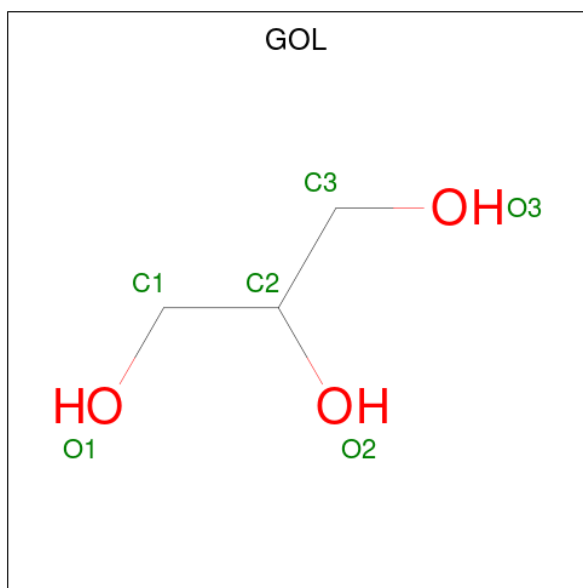
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Chain	Residue	Modelled	Actual	Comment	Reference
D	22	GLY	-	expression tag	UNP P0A903
D	23	SER	-	expression tag	UNP P0A903
D	24	HIS	-	expression tag	UNP P0A903
D	25	MET	-	expression tag	UNP P0A903

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total K 1 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Cl 1 1	0	0

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total O P 5 4 1	0	0

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	D	1	Total Na 1 1	0	0

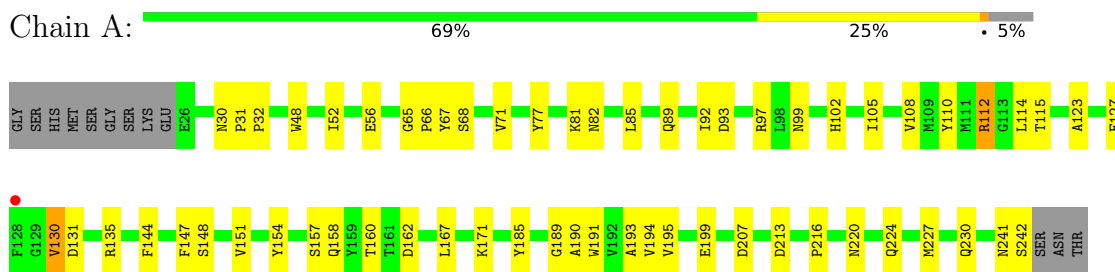
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	34	Total O 34 34	0	0
8	B	40	Total O 40 40	0	0
8	C	33	Total O 33 33	0	0
8	D	17	Total O 17 17	0	0

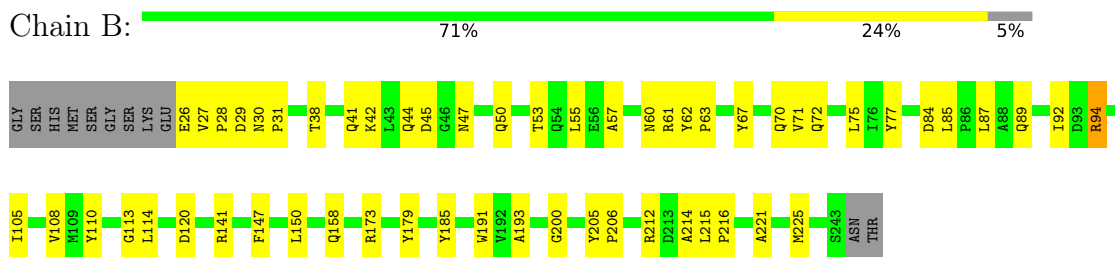
### 3 Residue-property plots [i](#)

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.

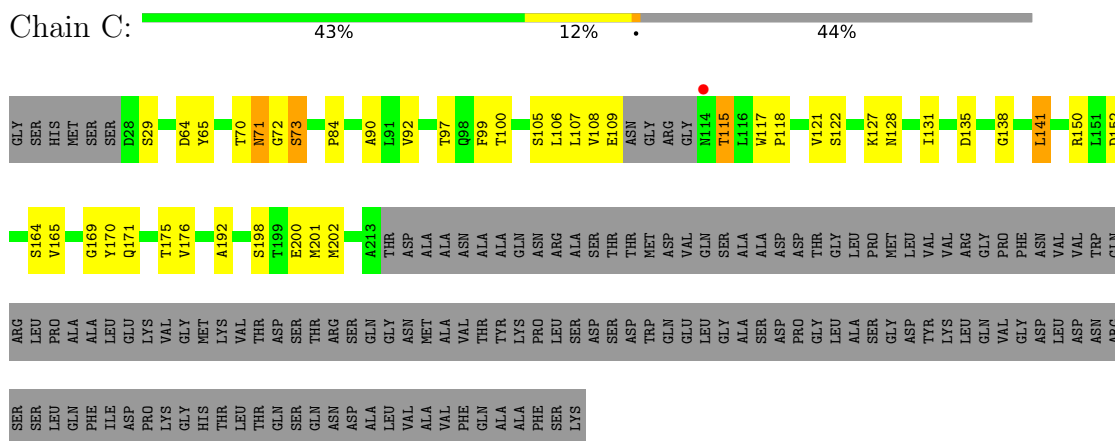
- Molecule 1: UPF0169 lipoprotein yfjO



- Molecule 1: UPF0169 lipoprotein yfjO



- Molecule 2: Lipoprotein 34



- Molecule 2: Lipoprotein 34



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.78Å 133.38Å 144.99Å 90.00° 100.15° 90.00°	Depositor
Resolution (Å)	97.45 – 2.90 51.70 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (97.45-2.90) 99.6 (51.70-2.90)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.73 (at 2.91Å)	Xtrriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.177 , 0.243 0.177 , 0.240	Depositor DCC
$R_{free}$ test set	1539 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.2	Xtrriage
Anisotropy	0.231	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6466	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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, K, NA, CL, PO4

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	0.70	0/1794	0.52	0/2438
1	B	0.67	0/1801	0.52	0/2447
2	C	0.70	0/1410	0.55	0/1923
2	D	0.75	3/1439 (0.2%)	0.52	0/1962
All	All	0.70	3/6444 (0.0%)	0.53	0/8770

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	60	VAL	CB-CG1	-5.96	1.40	1.52
2	D	60	VAL	CB-CG2	-5.50	1.41	1.52
2	D	129	TYR	CE2-CZ	-5.13	1.31	1.38

There are no bond angle outliers.

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	1754	0	1694	44	0
1	B	1761	0	1699	53	0
2	C	1383	0	1355	49	0
2	D	1412	0	1385	71	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
4	D	12	0	16	2	0
5	B	1	0	0	0	0
6	C	5	0	0	0	0
7	D	1	0	0	0	0
8	A	34	0	0	1	0
8	B	40	0	0	3	0
8	C	33	0	0	0	0
8	D	17	0	0	1	0
All	All	6466	0	6165	202	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 16.

All (202) 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:B:191:TRP:HB2	1:B:225:MET:CE	1.62	1.28
2:D:70:THR:HA	2:D:71:ASN:HB2	1.16	1.13
1:B:191:TRP:CB	1:B:225:MET:HE1	1.80	1.12
1:B:191:TRP:HB2	1:B:225:MET:HE1	1.24	1.08
1:B:191:TRP:CB	1:B:225:MET:CE	2.33	1.03
1:A:105:ILE:HG22	2:D:57:ILE:HD11	1.42	0.99
1:A:92:ILE:HG23	1:A:108:VAL:HG13	1.49	0.94
2:D:170:TYR:N	2:D:171:GLN:HB2	1.83	0.94
1:A:130:VAL:O	1:A:130:VAL:HG23	1.71	0.89
1:B:191:TRP:HB2	1:B:225:MET:HE3	1.52	0.89
2:D:70:THR:HA	2:D:71:ASN:CB	2.03	0.89
2:C:105:SER:HA	2:C:202:MET:HE3	1.55	0.88
1:B:26:GLU:HG2	1:B:27:VAL:H	1.43	0.83
1:B:92:ILE:HG23	1:B:108:VAL:HG12	1.60	0.81
2:C:169:GLY:HA3	2:C:170:TYR:C	2.02	0.79
1:B:158:GLN:HE22	2:D:192:ALA:HB2	1.47	0.78
1:B:67:TYR:O	1:B:71:VAL:HG23	1.83	0.78
1:A:92:ILE:HG23	1:A:108:VAL:CG1	2.15	0.76
1:A:89:GLN:CG	1:A:115:THR:HG21	2.17	0.75
2:D:70:THR:CA	2:D:71:ASN:HB2	2.09	0.75
2:D:192:ALA:HB3	8:D:346:HOH:O	1.87	0.74
1:A:89:GLN:HG3	1:A:115:THR:HG21	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:70:THR:HB	2:D:72:GLY:N	2.03	0.73
1:B:70:GLN:HB2	2:D:87[A]:GLN:NE2	2.02	0.73
1:B:191:TRP:CB	1:B:225:MET:HE3	2.09	0.72
1:A:158:GLN:HE22	2:C:192:ALA:HB2	1.53	0.72
2:C:105:SER:CB	2:C:175:THR:HG22	2.19	0.72
1:B:41:GLN:O	1:B:45:ASP:HB2	1.90	0.71
2:D:53:PRO:HG2	2:D:56:MET:SD	2.31	0.71
2:C:115:THR:HG23	2:C:115:THR:O	1.89	0.71
1:B:84:ASP:HB3	1:B:87:LEU:HD12	1.74	0.70
1:B:77:TYR:CG	2:D:84:PRO:HD3	2.26	0.70
2:C:170:TYR:CG	2:C:171:GLN:HA	2.26	0.70
2:D:127:LYS:O	2:D:128:ASN:HB2	1.91	0.69
1:B:29:ASP:OD1	1:B:29:ASP:C	2.31	0.68
1:B:44:GLN:O	2:D:32:LYS:NZ	2.22	0.67
2:C:109:GLU:O	2:C:170:TYR:HE2	1.78	0.67
2:D:170:TYR:HA	2:D:171:GLN:O	1.94	0.67
2:C:117:TRP:O	2:C:121:VAL:HG23	1.96	0.66
2:D:170:TYR:HA	2:D:171:GLN:C	2.16	0.65
1:B:30:ASN:HB3	1:B:31:PRO:HD2	1.79	0.64
1:B:47:ASN:OD1	1:B:50:GLN:HB2	1.97	0.64
2:C:105:SER:HB3	2:C:175:THR:HG22	1.77	0.64
2:D:109:GLU:HB3	2:D:171:GLN:H	1.63	0.64
1:A:151:VAL:HG12	1:A:160:THR:HG23	1.80	0.64
2:C:115:THR:O	2:C:115:THR:CG2	2.46	0.64
1:B:191:TRP:HB3	1:B:225:MET:CE	2.25	0.64
1:B:158:GLN:NE2	2:D:192:ALA:HB2	2.14	0.63
2:D:168:GLN:C	2:D:170:TYR:H	2.00	0.62
2:D:117:TRP:N	2:D:118:PRO:HD2	2.15	0.62
1:A:227:MET:HB3	1:A:230:GLN:OE1	2.00	0.62
2:C:105:SER:HB2	2:C:175:THR:HG22	1.81	0.62
1:B:30:ASN:HB3	1:B:31:PRO:CD	2.29	0.61
2:C:70:THR:O	2:C:71:ASN:HB2	2.00	0.61
2:D:169:GLY:C	2:D:171:GLN:HB2	2.19	0.61
2:C:170:TYR:CD2	2:C:171:GLN:HA	2.36	0.61
1:A:67:TYR:O	1:A:71:VAL:HG23	2.01	0.61
1:A:213:ASP:O	1:A:216:PRO:HD2	2.00	0.61
1:B:27:VAL:HG11	1:B:57:ALA:HB2	1.82	0.61
1:A:32:PRO:HD2	8:A:256:HOH:O	2.00	0.60
2:D:176:VAL:HG23	2:D:202:MET:HE1	1.82	0.60
1:B:57:ALA:O	1:B:61:ARG:HB2	2.01	0.60
2:C:117:TRP:HE3	2:C:165:VAL:HG23	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:127:LYS:O	2:C:128:ASN:HB2	2.02	0.59
1:B:221:ALA:O	1:B:225:MET:HG2	2.02	0.59
1:B:113:GLY:HA3	1:B:147:PHE:CE1	2.38	0.59
1:A:220:ASN:O	1:A:224:GLN:HG2	2.02	0.58
2:C:97:THR:HG22	2:C:106:LEU:HD12	1.84	0.58
1:A:144:PHE:CZ	2:C:65:TYR:HA	2.38	0.58
2:C:176:VAL:HG23	2:C:202:MET:HE1	1.84	0.58
2:D:133:GLN:NE2	2:D:142:THR:OG1	2.36	0.58
1:B:179:TYR:HD1	1:B:214:ALA:HB2	1.68	0.58
1:A:189:GLY:HA2	1:A:191:TRP:CZ2	2.39	0.58
1:B:110:TYR:HB2	1:B:150:LEU:HD22	1.86	0.57
2:D:133:GLN:O	2:D:141:LEU:HA	2.04	0.57
1:A:207:ASP:OD2	1:A:207:ASP:N	2.32	0.57
2:D:114:ASN:HB3	2:D:170:TYR:OH	2.05	0.57
2:C:108:VAL:O	2:C:171:GLN:HB2	2.05	0.56
1:B:120:ASP:OD2	1:B:173:ARG:NH1	2.38	0.56
2:D:150:ARG:NH2	2:D:152:ASP:OD2	2.38	0.56
2:D:121:VAL:O	2:D:125:GLN:HG3	2.05	0.56
1:A:89:GLN:HE21	1:A:115:THR:HG22	1.71	0.56
2:D:171:GLN:NE2	2:D:171:GLN:HA	2.21	0.56
1:B:200:GLY:HA3	8:B:4:HOH:O	2.05	0.56
2:D:189:ALA:HA	2:D:194:MET:CE	2.37	0.55
2:C:109:GLU:O	2:C:170:TYR:CE2	2.57	0.55
1:B:26:GLU:CG	1:B:27:VAL:H	2.18	0.55
1:B:26:GLU:HG2	1:B:27:VAL:N	2.19	0.55
1:A:93:ASP:OD1	1:A:112:ARG:NH1	2.41	0.54
2:D:69:VAL:O	2:D:70:THR:HG23	2.08	0.54
2:C:64:ASP:O	2:C:65:TYR:HB2	2.09	0.53
2:C:169:GLY:HA3	2:C:171:GLN:N	2.23	0.52
1:A:81:LYS:HA	1:A:81:LYS:HE2	1.90	0.52
1:A:154:TYR:CE1	2:D:57:ILE:HD12	2.45	0.52
2:D:208:GLY:HA2	2:D:211:LYS:HE2	1.90	0.52
2:C:107:LEU:HD12	2:C:171:GLN:HE21	1.75	0.52
2:D:189:ALA:HA	2:D:194:MET:HE3	1.92	0.52
1:A:110:TYR:OH	1:A:162:ASP:OD2	2.23	0.52
1:B:191:TRP:HB3	1:B:225:MET:HE1	1.81	0.51
1:B:191:TRP:C	1:B:225:MET:HE1	2.30	0.51
2:C:135:ASP:OD2	2:C:138:GLY:HA3	2.10	0.51
2:C:169:GLY:N	2:C:171:GLN:O	2.43	0.51
1:A:131:ASP:OD2	1:A:135:ARG:NH2	2.43	0.51
2:D:170:TYR:N	2:D:171:GLN:CB	2.67	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:LYS:HE3	2:C:65:TYR:HB3	1.92	0.51
1:A:52:ILE:O	1:A:56:GLU:HB2	2.11	0.51
2:C:72:GLY:HA3	2:C:73:SER:OG	2.11	0.51
2:D:176:VAL:HG23	2:D:202:MET:CE	2.41	0.50
2:D:170:TYR:C	2:D:170:TYR:CD1	2.85	0.50
1:A:48:TRP:CB	1:A:82:ASN:HB2	2.41	0.49
1:A:92:ILE:HG21	1:A:112:ARG:HB2	1.94	0.49
1:B:61:ARG:C	1:B:63:PRO:HD3	2.33	0.49
2:D:97:THR:O	2:D:98:GLN:HB3	2.13	0.49
1:B:85:LEU:O	1:B:89:GLN:HG3	2.12	0.49
2:C:117:TRP:HH2	2:C:141:LEU:HD23	1.76	0.49
1:A:144:PHE:CE2	2:C:65:TYR:HA	2.48	0.49
2:C:131:ILE:HG23	2:C:141:LEU:HD12	1.95	0.49
2:C:97:THR:HG22	2:C:106:LEU:CD1	2.42	0.49
2:D:70:THR:HB	2:D:72:GLY:H	1.76	0.49
2:D:109:GLU:HB3	2:D:171:GLN:N	2.28	0.49
1:B:29:ASP:OD1	1:B:29:ASP:O	2.30	0.49
2:D:171:GLN:NE2	2:D:171:GLN:CA	2.75	0.48
2:D:119:GLN:O	2:D:123:VAL:HG23	2.14	0.48
1:A:123:ALA:O	1:A:127:PHE:HD2	1.96	0.48
2:C:131:ILE:HG12	2:C:141:LEU:CD1	2.44	0.48
2:D:117:TRP:O	2:D:121:VAL:HG23	2.14	0.48
1:A:66:PRO:HG2	1:A:67:TYR:CD1	2.49	0.48
1:B:185:TYR:CD2	1:B:193:ALA:HB1	2.48	0.48
1:A:189:GLY:HA2	1:A:191:TRP:CH2	2.49	0.48
2:C:64:ASP:O	2:C:65:TYR:CB	2.62	0.47
2:C:117:TRP:CH2	2:C:141:LEU:CD2	2.97	0.47
2:C:150:ARG:HB3	2:C:152:ASP:OD1	2.14	0.47
2:D:117:TRP:N	2:D:118:PRO:CD	2.77	0.47
1:A:99:ASN:HB3	1:A:102:HIS:HB2	1.97	0.47
2:D:99:PHE:CZ	2:D:101:GLY:O	2.67	0.47
1:A:48:TRP:HB3	1:A:82:ASN:HB2	1.97	0.47
2:D:117:TRP:CH2	2:D:141:LEU:HD23	2.49	0.47
2:D:69:VAL:CG2	2:D:70:THR:N	2.78	0.47
2:D:169:GLY:C	2:D:171:GLN:CG	2.84	0.47
2:C:117:TRP:HH2	2:C:141:LEU:CD2	2.28	0.46
2:D:87[A]:GLN:NE2	2:D:87[A]:GLN:H	2.12	0.46
2:D:168:GLN:C	2:D:170:TYR:N	2.67	0.46
2:C:90:ALA:C	2:C:92:VAL:H	2.19	0.46
1:B:94:ARG:HA	1:B:94:ARG:HD3	1.70	0.46
2:D:171:GLN:HA	2:D:171:GLN:HE21	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:LEU:HB3	1:A:115:THR:HG23	1.98	0.46
1:A:185:TYR:CD2	1:A:193:ALA:HB1	2.50	0.46
1:A:190:ALA:O	1:A:194:VAL:HG23	2.15	0.46
2:C:117:TRP:HB3	2:C:118:PRO:HD3	1.99	0.45
2:D:69:VAL:HG22	2:D:70:THR:N	2.31	0.45
2:D:169:GLY:O	2:D:171:GLN:HG2	2.17	0.45
2:D:87[A]:GLN:H	2:D:87[A]:GLN:CD	2.19	0.45
1:B:205:TYR:N	1:B:206:PRO:CD	2.80	0.45
2:D:33:ARG:NH2	2:D:154:ASP:OD2	2.50	0.45
1:B:27:VAL:HG11	1:B:57:ALA:CB	2.47	0.44
2:D:89:LEU:HD13	2:D:203:ASN:ND2	2.32	0.44
1:A:68:SER:O	1:A:71:VAL:HB	2.17	0.44
2:C:117:TRP:CH2	2:C:141:LEU:HD23	2.52	0.44
2:D:89:LEU:HD13	2:D:203:ASN:HD21	1.83	0.44
1:A:241:ASN:O	1:A:242:SER:C	2.55	0.44
2:D:169:GLY:C	2:D:171:GLN:HG2	2.38	0.44
1:A:99:ASN:O	1:A:105:ILE:HD13	2.18	0.44
1:B:72:GLN:HG2	8:B:8:HOH:O	2.18	0.43
2:D:90:ALA:HB2	2:D:97:THR:HG23	1.99	0.43
1:A:195:VAL:O	1:A:199:GLU:HB2	2.19	0.43
2:D:102:ASP:OD1	2:D:102:ASP:N	2.48	0.43
2:D:145:TRP:CD1	2:D:158:ARG:NH1	2.86	0.43
2:C:107:LEU:HD22	2:C:107:LEU:N	2.34	0.43
2:D:169:GLY:C	2:D:171:GLN:CB	2.87	0.43
1:B:215:LEU:N	1:B:216:PRO:CD	2.80	0.43
1:A:77:TYR:CG	2:C:84:PRO:HD3	2.54	0.43
1:B:70:GLN:OE1	2:D:87[A]:GLN:NE2	2.39	0.43
1:A:30:ASN:HB3	1:A:31:PRO:CD	2.49	0.43
1:B:62:TYR:N	1:B:63:PRO:HD3	2.33	0.42
2:D:161:TYR:OH	2:D:197:TYR:O	2.30	0.42
2:D:38:ASP:H	4:D:2:GOL:C1	2.32	0.42
1:A:114:LEU:HD23	1:A:114:LEU:HA	1.76	0.42
2:C:117:TRP:CE3	2:C:165:VAL:HG23	2.53	0.42
2:D:52:ALA:HA	2:D:53:PRO:HD2	1.86	0.42
2:C:150:ARG:HH11	2:C:200:GLU:CD	2.22	0.42
1:B:212:ARG:NH2	2:D:48:ALA:O	2.50	0.42
2:C:72:GLY:HA3	2:C:73:SER:HA	1.72	0.42
1:B:105:ILE:HD12	1:B:105:ILE:HA	1.81	0.42
1:A:147:PHE:HB2	1:A:167:LEU:HD21	2.02	0.41
2:C:105:SER:CA	2:C:202:MET:HE3	2.38	0.41
2:C:169:GLY:CA	2:C:170:TYR:C	2.81	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:176:VAL:HG11	2:C:201:MET:HB3	2.02	0.41
1:B:27:VAL:HA	1:B:28:PRO:HD3	1.68	0.41
1:B:55:LEU:HB3	1:B:75:LEU:HG	2.03	0.41
1:B:110:TYR:CE2	1:B:114:LEU:HD11	2.55	0.41
2:C:117:TRP:CH2	2:C:141:LEU:HD22	2.55	0.41
2:D:99:PHE:O	2:D:100:THR:HG23	2.20	0.41
1:B:38:THR:O	1:B:42:LYS:HG2	2.19	0.41
2:C:99:PHE:HE1	2:C:198:SER:HB3	1.85	0.41
1:B:141:ARG:NH2	8:B:256:HOH:O	2.48	0.41
1:B:77:TYR:CD2	2:D:84:PRO:HD3	2.55	0.40
2:D:96:ARG:NH1	4:D:345:GOL:H11	2.36	0.40
1:A:89:GLN:HG2	1:A:115:THR:HG21	1.97	0.40
1:B:191:TRP:CA	1:B:225:MET:HE1	2.48	0.40
2:D:170:TYR:CA	2:D:171:GLN:HB2	2.51	0.40
2:D:178:LEU:HD21	2:D:181:LEU:HD22	2.03	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	215/229 (94%)	204 (95%)	9 (4%)	2 (1%)	17	48
1	B	216/229 (94%)	206 (95%)	10 (5%)	0	100	100
2	C	179/323 (55%)	161 (90%)	17 (10%)	1 (1%)	25	58
2	D	183/323 (57%)	162 (88%)	18 (10%)	3 (2%)	9	32
All	All	793/1104 (72%)	733 (92%)	54 (7%)	6 (1%)	19	51

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	170	TYR
2	C	71	ASN
2	D	128	ASN
2	D	61	THR
1	A	65	GLY
1	A	130	VAL

### 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	182/192 (95%)	178 (98%)	4 (2%)	52	81
1	B	183/192 (95%)	180 (98%)	3 (2%)	62	86
2	C	145/260 (56%)	138 (95%)	7 (5%)	25	58
2	D	148/260 (57%)	140 (95%)	8 (5%)	22	54
All	All	658/904 (73%)	636 (97%)	22 (3%)	37	72

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	ARG
1	A	112	ARG
1	A	148	SER
1	A	157	SER
1	B	53	THR
1	B	60	ASN
1	B	94	ARG
2	C	29	SER
2	C	73	SER
2	C	100	THR
2	C	115	THR
2	C	122	SER
2	C	141	LEU
2	C	164	SER
2	D	64	ASP
2	D	103	THR

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Mol	Chain	Res	Type
2	D	108	VAL
2	D	116	LEU
2	D	140	THR
2	D	151	LEU
2	D	162	GLN
2	D	193	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	158	GLN
1	B	158	GLN
2	C	119	GLN
2	D	114	ASN
2	D	133	GLN
2	D	171	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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
4	GOL	D	2	-	5,5,5	0.35	0	5,5,5	0.36	0
4	GOL	A	4	-	5,5,5	0.36	0	5,5,5	0.19	0
4	GOL	D	345	-	5,5,5	0.43	0	5,5,5	0.20	0
4	GOL	B	3	-	5,5,5	0.36	0	5,5,5	0.34	0
6	PO4	C	1	-	4,4,4	0.89	0	6,6,6	0.33	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
4	GOL	A	4	-	-	3/4/4/4	-
4	GOL	D	345	-	-	2/4/4/4	-
4	GOL	B	3	-	-	2/4/4/4	-
4	GOL	D	2	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	345	GOL	O1-C1-C2-C3
4	D	2	GOL	O1-C1-C2-C3
4	A	4	GOL	C1-C2-C3-O3
4	B	3	GOL	O1-C1-C2-C3
4	D	2	GOL	C1-C2-C3-O3
4	B	3	GOL	O1-C1-C2-O2
4	D	345	GOL	O1-C1-C2-O2
4	D	2	GOL	O1-C1-C2-O2
4	A	4	GOL	O2-C2-C3-O3
4	D	2	GOL	O2-C2-C3-O3
4	A	4	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	2	GOL	1	0
4	D	345	GOL	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 [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	217/229 (94%)	-0.46	1 (0%) 91   91	17, 31, 53, 70	0
1	B	218/229 (95%)	-0.54	0 100   100	15, 29, 58, 73	0
2	C	182/323 (56%)	-0.45	1 (0%) 91   91	17, 36, 63, 72	0
2	D	186/323 (57%)	-0.03	3 (1%) 72   71	17, 44, 76, 93	0
All	All	803/1104 (72%)	-0.38	5 (0%) 89   89	15, 34, 64, 93	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	114	ASN	3.1
2	C	114	ASN	2.6
2	D	170	TYR	2.6
2	D	169	GLY	2.3
1	A	128	PHE	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 monosaccharides 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
6	PO4	C	1	5/5	0.81	0.17	91,92,93,93	0
4	GOL	B	3	6/6	0.83	0.24	77,78,79,79	0
4	GOL	A	4	6/6	0.83	0.23	81,87,88,90	0
5	CL	B	1	1/1	0.90	0.26	40,40,40,40	0
4	GOL	D	345	6/6	0.91	0.19	58,61,63,64	0
4	GOL	D	2	6/6	0.91	0.24	56,63,66,67	0
3	K	A	1	1/1	0.93	0.10	61,61,61,61	0
7	NA	D	1	1/1	0.97	0.15	13,13,13,13	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.