



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:00 PM GMT

PDB ID : 1RLM
Title : Crystal Structure of ybiV from Escherichia coli K12
Authors : Roberts, A.; Lee, S.Y.; McCullagh, E.; Silversmith, R.E.; Wemmer, D.E.
Deposited on : 2003-11-26
Resolution : 1.90 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

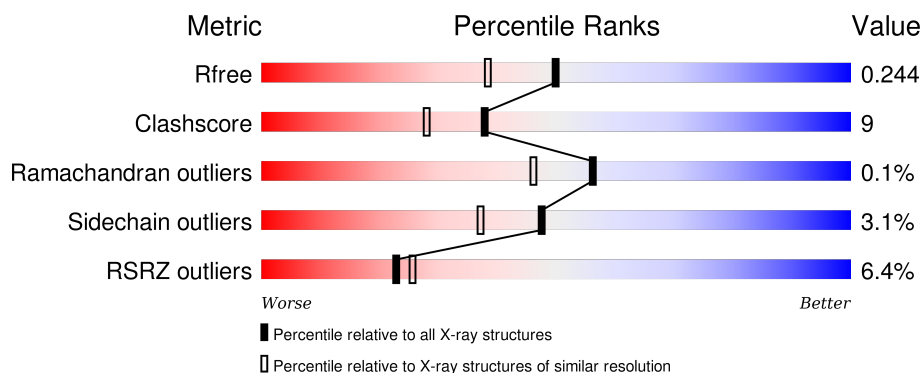
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	271	<div> <div>5%</div> <div>78%</div> <div>21%</div> <div>.</div> </div>
1	B	271	<div> <div>8%</div> <div>82%</div> <div>16%</div> <div>..</div> </div>
1	C	271	<div> <div>11%</div> <div>75%</div> <div>23%</div> <div>..</div> </div>
1	D	271	<div> <div>2%</div> <div>81%</div> <div>16%</div> <div>..</div> </div>

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
3	GOL	A	804	-	X	-	X
3	GOL	A	806	-	X	-	X
3	GOL	B	807	-	X	-	X
3	GOL	C	808	-	X	-	-
3	GOL	C	809	-	X	-	X
3	GOL	D	810	-	X	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9154 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 Phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2134	1369	362	395	8			
1	B	269	Total	C	N	O	S	0	0	0
			2134	1369	362	395	8			
1	C	269	Total	C	N	O	S	0	0	0
			2134	1369	362	395	8			
1	D	269	Total	C	N	O	S	0	0	0
			2134	1369	362	395	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	SER	ENGINEERED	UNP P75792
A	267	TYR	SER	ENGINEERED	UNP P75792
B	2	ALA	SER	ENGINEERED	UNP P75792
B	267	TYR	SER	ENGINEERED	UNP P75792
C	2	ALA	SER	ENGINEERED	UNP P75792
C	267	TYR	SER	ENGINEERED	UNP P75792
D	2	ALA	SER	ENGINEERED	UNP P75792
D	267	TYR	SER	ENGINEERED	UNP P75792

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

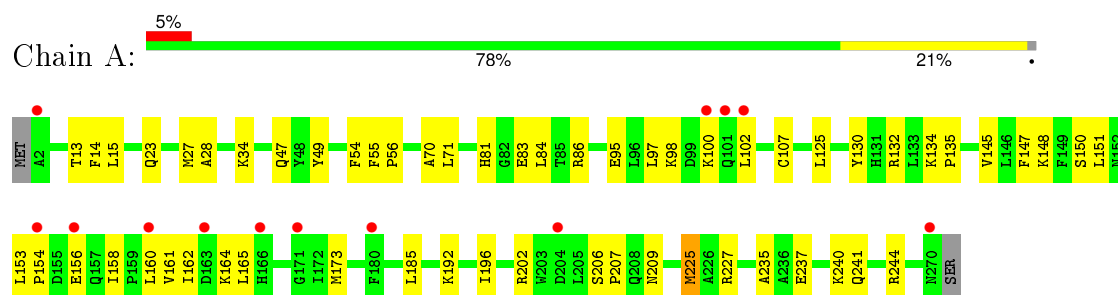
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	151	Total	O	0	0
			151	151		
4	B	150	Total	O	0	0
			150	150		
4	C	107	Total	O	0	0
			107	107		
4	D	170	Total	O	0	0
			170	170		

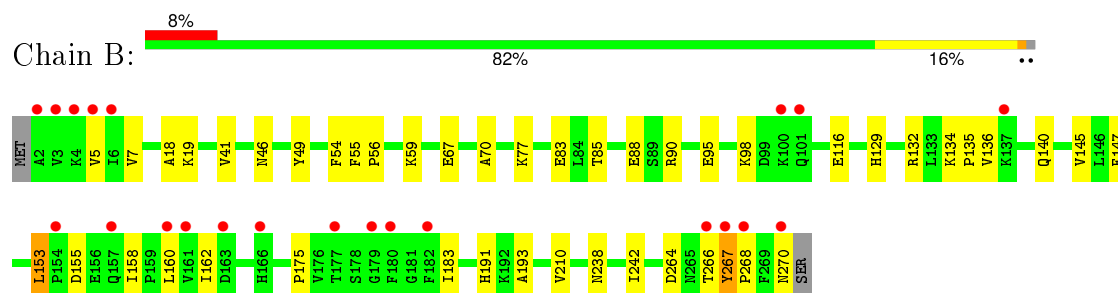
3 Residue-property plots [i](#)

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.

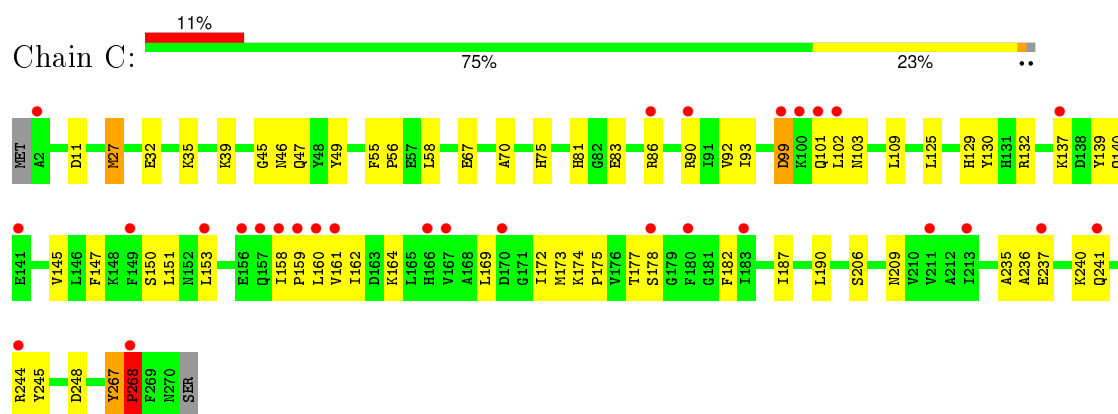
• Molecule 1: Phosphatase



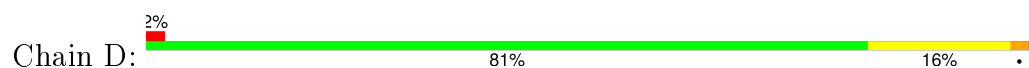
• Molecule 1: Phosphatase

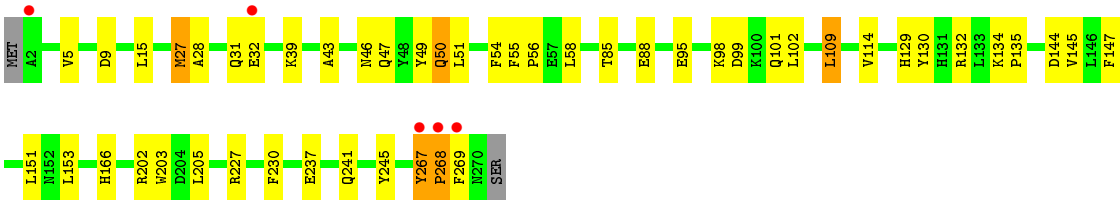


• Molecule 1: Phosphatase



• Molecule 1: Phosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.82Å 91.61Å 186.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 1.90 19.96 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.1 (19.96-1.90) 97.2 (19.96-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.10 (at 1.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.248 0.217 , 0.244	Depositor DCC
R_{free} test set	9467 reflections (9.98%)	DCC
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 96269 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9154	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: GOL, MG

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	0.34	0/2179	0.60	0/2946
1	B	0.37	0/2179	0.67	2/2946 (0.1%)
1	C	0.34	0/2179	0.84	5/2946 (0.2%)
1	D	0.39	0/2179	0.71	2/2946 (0.1%)
All	All	0.36	0/8716	0.71	9/11784 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	267	TYR	C-N-CD	-19.40	77.92	120.60
1	C	268	PRO	CA-N-CD	-17.61	86.84	111.50
1	D	267	TYR	C-N-CD	-13.11	91.75	120.60
1	C	268	PRO	N-CA-CB	12.14	117.87	103.30
1	C	268	PRO	N-CD-CG	11.79	120.89	103.20
1	B	267	TYR	C-N-CD	-11.70	94.86	120.60
1	C	267	TYR	C-N-CA	9.24	160.81	122.00
1	D	267	TYR	C-N-CA	9.08	160.15	122.00
1	B	267	TYR	C-N-CA	7.13	151.94	122.00

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	2134	0	2124	38	0
1	B	2134	0	2124	33	0
1	C	2134	0	2124	47	0
1	D	2134	0	2124	39	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	12	0	8	1	0
3	B	6	0	4	0	0
3	C	12	0	8	3	0
3	D	6	0	4	0	0
4	A	151	0	0	4	0
4	B	150	0	0	1	0
4	C	107	0	0	2	0
4	D	170	0	0	3	0
All	All	9154	0	8520	157	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 9.

All (157) 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:99:ASP:OD1	1:D:101:GLN:HG2	1.76	0.85
1:A:102:LEU:HD11	1:A:151:LEU:HD23	1.57	0.83
1:C:83:GLU:HG3	1:C:145:VAL:HG13	1.59	0.82
1:B:267:TYR:CD2	1:B:268:PRO:HD3	2.20	0.76
1:A:13:THR:HG22	1:A:14:PHE:N	1.98	0.76
1:C:267:TYR:CD2	1:C:268:PRO:HD3	2.23	0.73
1:C:102:LEU:HD23	1:C:103:ASN:N	2.05	0.71
1:C:102:LEU:HD11	1:C:151:LEU:HD23	1.71	0.71
1:D:27:MET:HE3	1:D:58:LEU:HD11	1.74	0.70
1:C:11:ASP:HB2	3:C:809:GOL:H11	1.73	0.69
1:C:102:LEU:HD21	1:C:151:LEU:HG	1.74	0.67
1:A:164:LYS:HG2	4:A:903:HOH:O	1.93	0.66
1:A:156:GLU:CD	1:A:156:GLU:H	2.00	0.65
1:A:227:ARG:HD2	1:A:244:ARG:NH2	2.12	0.64
1:A:207:PRO:HB3	1:A:225:MET:HE2	1.80	0.63
1:C:161:VAL:HA	1:C:164:LYS:HE2	1.81	0.63
1:A:154:PRO:HB3	1:A:156:GLU:OE2	1.99	0.62
1:C:178:SER:HB3	1:C:182:PHE:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LEU:HD12	3:A:806:GOL:O2	2.01	0.60
1:B:85:THR:OG1	1:B:88:GLU:HG3	2.02	0.59
1:D:49:TYR:CG	1:D:132:ARG:HD3	2.38	0.58
1:B:162:ILE:HG23	1:B:175:PRO:HB2	1.85	0.58
1:C:187:ILE:HB	1:C:190:LEU:HD12	1.86	0.57
1:C:90:ARG:HD3	1:C:140:GLN:NE2	2.19	0.57
1:B:264:ASP:HB2	1:B:266:THR:HG23	1.88	0.55
1:D:27:MET:CE	1:D:31:GLN:HE21	2.18	0.55
1:C:45:GLY:H	3:C:809:GOL:C3	2.18	0.55
1:A:153:LEU:HD23	1:A:153:LEU:N	2.21	0.55
1:A:165:LEU:HD13	1:A:185:LEU:HD21	1.88	0.55
1:D:50:GLN:HG2	1:D:129:HIS:O	2.06	0.55
1:A:237:GLU:O	1:A:241:GLN:HG3	2.07	0.55
1:B:55:PHE:N	1:B:56:PRO:HD3	2.23	0.54
1:A:49:TYR:CG	1:A:132:ARG:HD3	2.42	0.54
1:A:100:LYS:HE2	4:A:938:HOH:O	2.07	0.54
1:B:83:GLU:HG3	1:B:145:VAL:HG13	1.89	0.54
1:D:114:VAL:O	1:D:135:PRO:HA	2.07	0.54
1:D:153:LEU:HD23	1:D:153:LEU:N	2.23	0.54
1:A:227:ARG:HD2	1:A:244:ARG:HH22	1.71	0.54
1:C:86:ARG:HH11	1:C:86:ARG:HG2	1.72	0.53
1:C:160:LEU:HD23	4:C:871:HOH:O	2.08	0.53
1:A:55:PHE:N	1:A:56:PRO:HD3	2.24	0.53
1:D:166:HIS:HD2	4:D:912:HOH:O	1.92	0.52
1:B:49:TYR:CG	1:B:132:ARG:HD3	2.45	0.52
1:C:162:ILE:HD11	1:C:177:THR:HG22	1.92	0.51
1:D:5:VAL:HG23	1:D:205:LEU:HD23	1.93	0.51
1:C:153:LEU:N	1:C:153:LEU:HD23	2.26	0.51
1:D:99:ASP:HB3	1:D:102:LEU:HD22	1.93	0.51
1:C:27:MET:HE1	1:C:58:LEU:HD21	1.91	0.51
1:D:28:ALA:O	1:D:32:GLU:HG3	2.11	0.51
1:D:27:MET:CE	1:D:58:LEU:HD11	2.40	0.50
1:A:151:LEU:N	1:A:151:LEU:HD12	2.26	0.50
1:A:95:GLU:OE2	1:A:98:LYS:HD2	2.12	0.50
1:C:102:LEU:HD21	1:C:151:LEU:CG	2.41	0.50
1:C:237:GLU:O	1:C:241:GLN:HG3	2.11	0.50
1:B:267:TYR:CA	1:B:270:ASN:ND2	2.75	0.50
1:B:267:TYR:HA	1:B:270:ASN:ND2	2.27	0.50
1:C:49:TYR:CG	1:C:132:ARG:HD3	2.47	0.50
1:B:153:LEU:N	1:B:153:LEU:HD23	2.26	0.49
1:B:191:HIS:CE1	1:B:193:ALA:HB3	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:LYS:O	1:A:196:ILE:HG13	2.12	0.48
1:A:107:CYS:SG	1:A:148:LYS:HE2	2.53	0.48
1:A:83:GLU:HG3	1:A:145:VAL:HG13	1.94	0.48
1:D:132:ARG:HG2	4:D:812:HOH:O	2.14	0.48
1:C:93:ILE:HD12	1:C:139:TYR:HB3	1.95	0.48
1:A:153:LEU:CD1	1:A:161:VAL:HG21	2.44	0.47
1:B:18:ALA:O	1:B:19:LYS:HB2	2.13	0.47
1:D:227:ARG:O	1:D:227:ARG:HD2	2.14	0.47
1:A:206:SER:H	1:A:209:ASN:HD22	1.61	0.47
1:B:191:HIS:HB2	4:B:913:HOH:O	2.13	0.47
1:D:227:ARG:HG2	1:D:227:ARG:HH11	1.79	0.47
1:C:47:GLN:HB2	1:C:130:TYR:CD2	2.49	0.47
1:B:158:ILE:O	1:B:162:ILE:HG13	2.15	0.46
1:D:203:TRP:O	1:D:205:LEU:HD13	2.14	0.46
1:C:178:SER:HB3	1:C:182:PHE:CB	2.44	0.46
1:D:46:ASN:HB3	1:D:50:GLN:CG	2.45	0.46
1:B:238:ASN:O	1:B:242:ILE:HG13	2.15	0.46
1:B:160:LEU:HD23	1:B:160:LEU:C	2.36	0.46
1:C:150:SER:C	1:C:151:LEU:HD12	2.35	0.46
1:D:46:ASN:HB3	1:D:50:GLN:HG2	1.98	0.46
1:A:84:LEU:HD11	1:A:173:MET:CE	2.45	0.46
1:C:125:LEU:HD11	1:C:182:PHE:CZ	2.51	0.46
1:B:54:PHE:C	1:B:56:PRO:HD3	2.36	0.46
1:A:235:ALA:O	1:A:240:LYS:HE3	2.15	0.46
1:D:9:ASP:HA	1:D:43:ALA:O	2.16	0.45
1:B:160:LEU:HD23	1:B:160:LEU:O	2.16	0.45
1:C:11:ASP:CB	3:C:809:GOL:H11	2.43	0.45
1:B:267:TYR:CE2	1:B:268:PRO:HD3	2.51	0.45
1:C:162:ILE:HG23	1:C:175:PRO:HB2	1.99	0.45
1:D:245:TYR:CD2	1:D:268:PRO:HG3	2.52	0.45
1:C:132:ARG:HG2	4:C:836:HOH:O	2.16	0.44
1:C:235:ALA:O	1:C:240:LYS:HE3	2.18	0.44
1:B:95:GLU:OE2	1:B:98:LYS:HD2	2.18	0.44
1:D:46:ASN:OD1	1:D:129:HIS:HD2	2.00	0.44
1:B:5:VAL:HB	1:B:210:VAL:HG22	1.99	0.44
1:B:46:ASN:OD1	1:B:129:HIS:ND1	2.42	0.44
1:C:244:ARG:HG2	1:C:244:ARG:HH11	1.83	0.44
1:D:15:LEU:HD21	1:D:54:PHE:CE1	2.53	0.43
1:C:206:SER:H	1:C:209:ASN:HD22	1.66	0.43
1:C:92:VAL:HG22	1:C:169:LEU:HD22	2.01	0.43
1:D:134:LYS:HA	1:D:135:PRO:HD3	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:LEU:H	1:B:153:LEU:HD23	1.82	0.43
1:D:230:PHE:CZ	1:D:268:PRO:HG2	2.53	0.43
1:C:125:LEU:C	1:C:125:LEU:HD13	2.39	0.43
1:A:23:GLN:O	1:A:27:MET:HG2	2.18	0.43
1:A:34:LYS:HA	4:A:930:HOH:O	2.18	0.43
1:C:35:LYS:HE2	1:C:35:LYS:HB3	1.81	0.43
1:D:267:TYR:C	1:D:267:TYR:CD2	2.92	0.43
1:B:267:TYR:HB2	1:B:270:ASN:ND2	2.33	0.43
1:D:27:MET:HE1	1:D:58:LEU:HD21	2.00	0.43
1:C:32:GLU:HA	1:C:35:LYS:HE2	2.01	0.43
1:B:267:TYR:HB2	1:B:270:ASN:HD21	1.82	0.42
1:D:55:PHE:N	1:D:56:PRO:HD3	2.34	0.42
1:D:237:GLU:O	1:D:241:GLN:HG3	2.19	0.42
1:C:245:TYR:CE2	1:C:268:PRO:HG3	2.54	0.42
1:D:27:MET:HE2	1:D:31:GLN:HE21	1.85	0.42
1:C:70:ALA:O	1:C:81:HIS:HA	2.20	0.42
1:B:7:VAL:HG22	1:B:41:VAL:HB	2.01	0.42
1:A:49:TYR:CD1	1:A:132:ARG:HD3	2.54	0.42
1:D:144:ASP:OD2	1:D:145:VAL:N	2.52	0.42
1:D:28:ALA:HB3	4:D:897:HOH:O	2.20	0.42
1:D:5:VAL:CG2	1:D:205:LEU:HD23	2.50	0.42
1:C:125:LEU:HD11	1:C:182:PHE:HZ	1.84	0.41
1:C:55:PHE:N	1:C:56:PRO:HD3	2.34	0.41
1:A:154:PRO:CB	1:A:156:GLU:OE2	2.66	0.41
1:A:28:ALA:HB3	4:A:872:HOH:O	2.20	0.41
1:A:150:SER:C	1:A:151:LEU:HD12	2.40	0.41
1:B:155:ASP:HA	1:B:158:ILE:HG13	2.01	0.41
1:A:54:PHE:C	1:A:56:PRO:HD3	2.41	0.41
1:D:47:GLN:HB2	1:D:130:TYR:CD2	2.55	0.41
1:D:95:GLU:HA	1:D:98:LYS:HD3	2.03	0.41
1:C:158:ILE:HB	1:C:159:PRO:HD3	2.02	0.41
1:A:47:GLN:HB2	1:A:130:TYR:CD2	2.55	0.41
1:D:39:LYS:HD3	1:D:203:TRP:CE2	2.55	0.41
1:C:46:ASN:OD1	1:C:129:HIS:HD2	2.04	0.41
1:C:245:TYR:CD2	1:C:268:PRO:HG3	2.55	0.41
1:A:70:ALA:O	1:A:81:HIS:HA	2.20	0.41
1:C:99:ASP:C	1:C:101:GLN:H	2.22	0.41
1:B:134:LYS:HA	1:B:135:PRO:HD3	1.92	0.41
1:D:227:ARG:NH1	1:D:227:ARG:HG2	2.36	0.41
1:C:236:ALA:O	1:C:240:LYS:HG3	2.21	0.41
1:B:116:GLU:HG2	1:B:136:VAL:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LEU:HD21	1:A:151:LEU:CD2	2.50	0.41
1:B:158:ILE:HG23	1:B:183:ILE:HD11	2.03	0.41
1:A:13:THR:HG22	1:A:14:PHE:H	1.79	0.40
1:B:153:LEU:N	1:B:153:LEU:CD2	2.84	0.40
1:C:67:GLU:HB2	1:C:70:ALA:HB3	2.01	0.40
1:A:134:LYS:HA	1:A:135:PRO:HD3	1.88	0.40
1:C:172:ILE:HG22	1:C:173:MET:HG3	2.02	0.40
1:A:160:LEU:C	1:A:160:LEU:HD13	2.42	0.40
1:A:158:ILE:O	1:A:162:ILE:HG13	2.21	0.40
1:D:85:THR:OG1	1:D:88:GLU:HG3	2.22	0.40
1:C:86:ARG:NH1	1:C:86:ARG:HG2	2.35	0.40
1:B:67:GLU:HB2	1:B:70:ALA:HB3	2.04	0.40
1:D:50:GLN:NE2	1:D:51:LEU:HD23	2.36	0.40
1:C:174:LYS:HA	1:C:175:PRO:HD3	1.86	0.40
1:D:109:LEU:HB2	1:D:145:VAL:HB	2.03	0.40
1:B:90:ARG:HA	1:B:140:GLN:OE1	2.22	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	267/271 (98%)	255 (96%)	12 (4%)	0	100	100
1	B	267/271 (98%)	254 (95%)	13 (5%)	0	100	100
1	C	267/271 (98%)	254 (95%)	12 (4%)	1 (0%)	39	27
1	D	267/271 (98%)	255 (96%)	12 (4%)	0	100	100
All	All	1068/1084 (98%)	1018 (95%)	49 (5%)	1 (0%)	56	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	268	PRO

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	228/230 (99%)	221 (97%)	7 (3%)	47	37
1	B	228/230 (99%)	224 (98%)	4 (2%)	66	61
1	C	228/230 (99%)	219 (96%)	9 (4%)	39	27
1	D	228/230 (99%)	220 (96%)	8 (4%)	43	31
All	All	912/920 (99%)	884 (97%)	28 (3%)	47	37

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	71	LEU
1	A	86	ARG
1	A	97	LEU
1	A	147	PHE
1	A	202	ARG
1	A	225	MET
1	B	59	LYS
1	B	77	LYS
1	B	147	PHE
1	B	153	LEU
1	C	27	MET
1	C	39	LYS
1	C	75	HIS
1	C	99	ASP
1	C	109	LEU
1	C	137	LYS
1	C	147	PHE
1	C	248	ASP
1	C	268	PRO
1	D	27	MET

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Mol	Chain	Res	Type
1	D	50	GLN
1	D	109	LEU
1	D	147	PHE
1	D	151	LEU
1	D	202	ARG
1	D	268	PRO
1	D	269	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	HIS
1	A	101	GLN
1	A	110	GLN
1	A	208	GLN
1	A	209	ASN
1	B	78	GLN
1	B	87	HIS
1	B	270	ASN
1	C	129	HIS
1	C	208	GLN
1	C	209	ASN
1	D	31	GLN
1	D	50	GLN
1	D	78	GLN
1	D	101	GLN
1	D	129	HIS
1	D	152	ASN
1	D	166	HIS
1	D	209	ASN
1	D	241	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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
3	GOL	A	804	-	5,5,5	4.75	5 (100%)	5,5,5	5.64	3 (60%)
3	GOL	A	806	-	5,5,5	4.89	5 (100%)	5,5,5	5.58	3 (60%)
3	GOL	B	807	-	5,5,5	4.79	5 (100%)	5,5,5	5.69	3 (60%)
3	GOL	C	808	-	5,5,5	4.80	5 (100%)	5,5,5	5.69	3 (60%)
3	GOL	C	809	-	5,5,5	4.85	5 (100%)	5,5,5	5.67	3 (60%)
3	GOL	D	810	-	5,5,5	4.92	5 (100%)	5,5,5	5.64	3 (60%)

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
3	GOL	A	804	-	-	0/4/4/4	0/0/0/0
3	GOL	A	806	-	-	0/4/4/4	0/0/0/0
3	GOL	B	807	-	-	0/4/4/4	0/0/0/0
3	GOL	C	808	-	-	0/4/4/4	0/0/0/0
3	GOL	C	809	-	-	0/4/4/4	0/0/0/0
3	GOL	D	810	-	-	0/4/4/4	0/0/0/0

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	810	GOL	C3-C2	-8.51	1.19	1.52
3	C	809	GOL	C3-C2	-8.46	1.20	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	806	GOL	C3-C2	-8.40	1.20	1.52
3	B	807	GOL	C3-C2	-8.28	1.20	1.52
3	C	808	GOL	C3-C2	-8.26	1.20	1.52
3	A	804	GOL	C3-C2	-8.21	1.20	1.52
3	A	806	GOL	C1-C2	-3.61	1.38	1.52
3	C	809	GOL	C1-C2	-3.48	1.39	1.52
3	D	810	GOL	O2-C2	-3.42	1.33	1.43
3	D	810	GOL	C1-C2	-3.14	1.40	1.52
3	C	808	GOL	C1-C2	-3.14	1.40	1.52
3	B	807	GOL	C1-C2	-3.12	1.40	1.52
3	C	809	GOL	O2-C2	-3.02	1.34	1.43
3	A	804	GOL	C1-C2	-2.97	1.40	1.52
3	A	806	GOL	O2-C2	-2.97	1.34	1.43
3	C	808	GOL	O2-C2	-2.86	1.34	1.43
3	B	807	GOL	O2-C2	-2.83	1.35	1.43
3	A	804	GOL	O2-C2	-2.62	1.35	1.43
3	D	810	GOL	O3-C3	2.63	1.53	1.42
3	C	809	GOL	O3-C3	2.91	1.54	1.42
3	B	807	GOL	O3-C3	3.01	1.55	1.42
3	A	804	GOL	O3-C3	3.03	1.55	1.42
3	A	806	GOL	O3-C3	3.04	1.55	1.42
3	C	808	GOL	O3-C3	3.10	1.55	1.42
3	C	809	GOL	O1-C1	4.07	1.59	1.42
3	A	806	GOL	O1-C1	4.22	1.60	1.42
3	C	808	GOL	O1-C1	4.38	1.61	1.42
3	B	807	GOL	O1-C1	4.41	1.61	1.42
3	D	810	GOL	O1-C1	4.51	1.61	1.42
3	A	804	GOL	O1-C1	4.53	1.61	1.42

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	806	GOL	O1-C1-C2	2.92	124.33	110.18
3	D	810	GOL	O1-C1-C2	3.12	125.31	110.18
3	C	808	GOL	O1-C1-C2	3.13	125.37	110.18
3	B	807	GOL	O1-C1-C2	3.18	125.63	110.18
3	C	809	GOL	O1-C1-C2	3.31	126.26	110.18
3	A	804	GOL	O1-C1-C2	3.32	126.28	110.18
3	D	810	GOL	O2-C2-C3	6.37	137.86	108.65
3	C	808	GOL	O2-C2-C3	6.56	138.72	108.65
3	A	804	GOL	O2-C2-C3	6.56	138.75	108.65
3	B	807	GOL	O2-C2-C3	6.58	138.82	108.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	809	GOL	O2-C2-C3	6.71	139.44	108.65
3	A	806	GOL	O2-C2-C3	6.85	140.08	108.65
3	A	806	GOL	O3-C3-C2	9.98	158.57	110.18
3	C	809	GOL	O3-C3-C2	10.19	159.60	110.18
3	A	804	GOL	O3-C3-C2	10.20	159.67	110.18
3	B	807	GOL	O3-C3-C2	10.39	160.55	110.18
3	D	810	GOL	O3-C3-C2	10.39	160.57	110.18
3	C	808	GOL	O3-C3-C2	10.41	160.67	110.18

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	806	GOL	1	0
3	C	809	GOL	3	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	269/271 (99%)	0.49	13 (4%) 34 37	13, 24, 52, 83	0
1	B	269/271 (99%)	0.52	22 (8%) 14 16	12, 23, 61, 131	0
1	C	269/271 (99%)	0.69	29 (10%) 8 8	17, 30, 64, 118	0
1	D	269/271 (99%)	0.16	5 (1%) 70 73	12, 22, 43, 76	0
All	All	1076/1084 (99%)	0.47	69 (6%) 23 25	12, 25, 58, 131	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	5	VAL	8.8
1	B	2	ALA	8.7
1	B	3	VAL	7.0
1	B	101	GLN	6.4
1	B	160	LEU	6.1
1	A	101	GLN	5.7
1	A	2	ALA	5.6
1	C	100	LYS	5.2
1	C	180	PHE	5.0
1	B	4	LYS	4.9
1	C	160	LEU	4.8
1	C	156	GLU	4.7
1	A	100	LYS	4.0
1	B	6	ILE	4.0
1	B	100	LYS	3.8
1	A	154	PRO	3.8
1	A	163	ASP	3.6
1	C	2	ALA	3.6
1	A	180	PHE	3.5
1	B	180	PHE	3.5
1	C	101	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	166	HIS	3.3
1	D	268	PRO	3.1
1	C	157	GLN	3.1
1	B	268	PRO	3.0
1	C	167	VAL	3.0
1	D	2	ALA	3.0
1	C	102	LEU	3.0
1	B	182	PHE	2.9
1	A	102	LEU	2.9
1	C	137	LYS	2.9
1	C	183	ILE	2.7
1	A	156	GLU	2.7
1	D	267	TYR	2.7
1	B	179	GLY	2.7
1	A	204	ASP	2.7
1	C	268	PRO	2.7
1	A	270	ASN	2.6
1	C	213	ILE	2.6
1	B	177	THR	2.6
1	B	154	PRO	2.6
1	C	161	VAL	2.6
1	C	99	ASP	2.5
1	B	166	HIS	2.5
1	C	86	ARG	2.4
1	B	270	ASN	2.4
1	B	266	THR	2.4
1	B	267	TYR	2.3
1	C	149	PHE	2.3
1	B	137	LYS	2.3
1	C	170	ASP	2.3
1	B	163	ASP	2.3
1	C	158	ILE	2.3
1	A	171	GLY	2.3
1	C	244	ARG	2.3
1	B	157	GLN	2.2
1	C	241	GLN	2.2
1	C	141	GLU	2.2
1	C	90	ARG	2.2
1	A	160	LEU	2.2
1	C	153	LEU	2.2
1	B	161	VAL	2.2
1	D	32	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	166	HIS	2.1
1	C	211	VAL	2.1
1	C	237	GLU	2.1
1	D	269	PHE	2.1
1	C	178	SER	2.1
1	C	159	PRO	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 carbohydrates 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. 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
3	GOL	A	806	6/6	0.78	0.47	12.15	28,31,32,36	0
3	GOL	D	810	6/6	0.74	0.27	5.07	24,30,32,32	0
3	GOL	C	809	6/6	0.86	0.29	4.70	29,30,32,34	0
3	GOL	B	807	6/6	0.71	0.27	3.12	30,33,35,38	0
3	GOL	A	804	6/6	0.73	0.25	2.76	22,29,29,29	0
3	GOL	C	808	6/6	0.81	0.19	1.53	25,28,29,31	0
2	MG	A	800	1/1	0.97	0.09	-0.71	18,18,18,18	0
2	MG	B	801	1/1	0.97	0.07	-1.18	20,20,20,20	0
2	MG	D	802	1/1	0.98	0.06	-1.70	18,18,18,18	0
2	MG	C	803	1/1	0.95	0.09	-1.87	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.