



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 09:40 PM EDT

PDB ID : 3WSV  
Title : Crystal structure of minor L-lactate dehydrogenase from *Enterococcus mundtii* in the ligands-unbound form  
Authors : Matoba, Y.; Sugiyama, M.  
Deposited on : 2014-03-27  
Resolution : 2.38 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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
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)	:	2.36.2

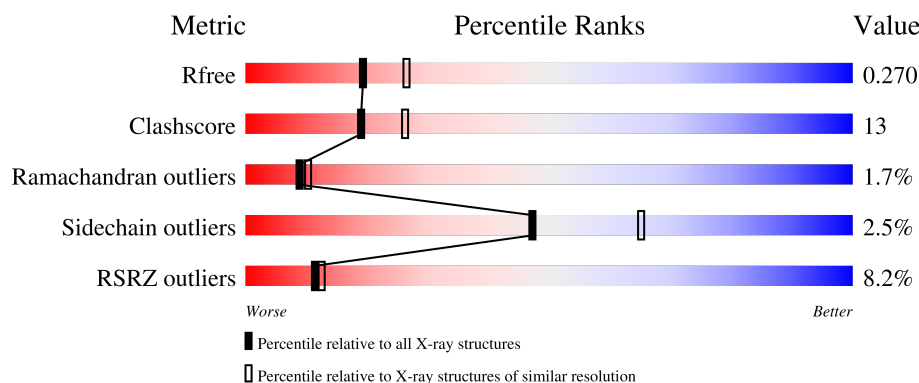
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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	322	<div> <div>11%</div> <div>65%</div> <div>31%</div> <div>..</div> </div>
1	B	322	<div> <div>5%</div> <div>74%</div> <div>23%</div> <div>..</div> </div>
1	C	322	<div> <div>10%</div> <div>67%</div> <div>27%</div> <div>..</div> </div>
1	D	322	<div> <div>6%</div> <div>78%</div> <div>20%</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
2	GOL	C	503	-	-	-	X

## 2 Entry composition [i](#)

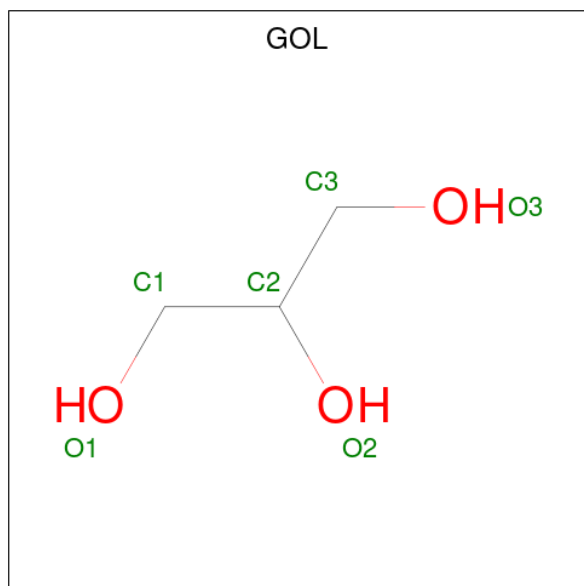
There are 3 unique types of molecules in this entry. The entry contains 10138 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-lactate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	0	0
			2400	1513	407	472	8			
1	B	317	Total	C	N	O	S	0	0	0
			2441	1538	415	481	7			
1	C	312	Total	C	N	O	S	0	0	0
			2400	1513	407	472	8			
1	D	317	Total	C	N	O	S	0	0	0
			2441	1538	415	481	7			

- Molecule 2 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
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

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

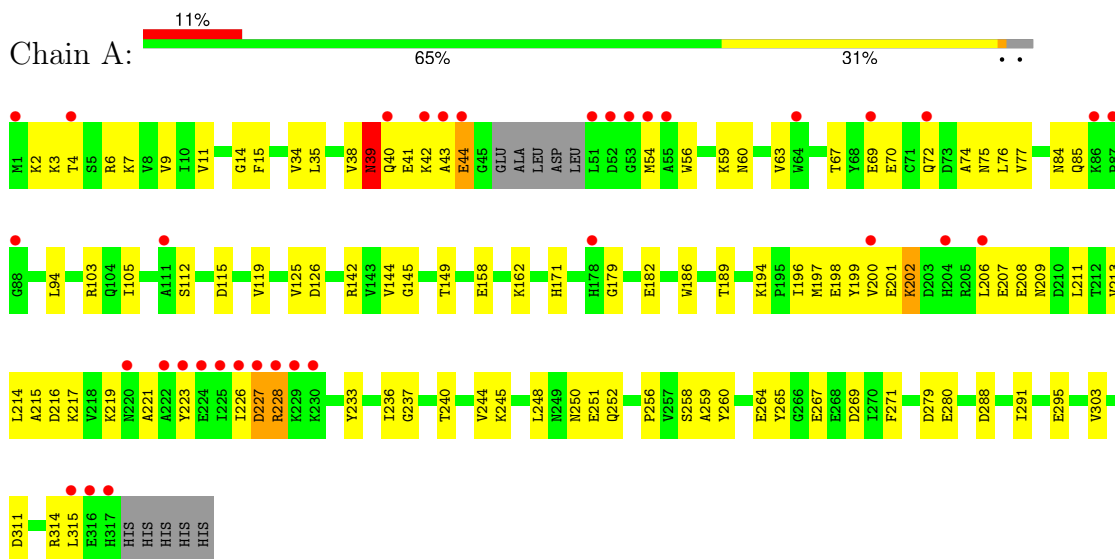
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	86	Total O 86 86	0	0
3	B	121	Total O 121 121	0	0
3	C	79	Total O 79 79	0	0
3	D	122	Total O 122 122	0	0

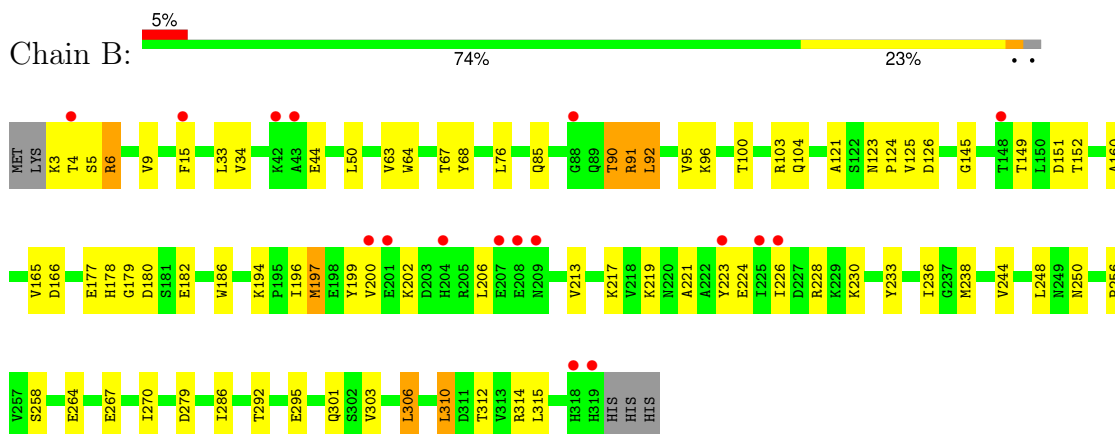
### 3 Residue-property plots [i](#)

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: L-lactate dehydrogenase

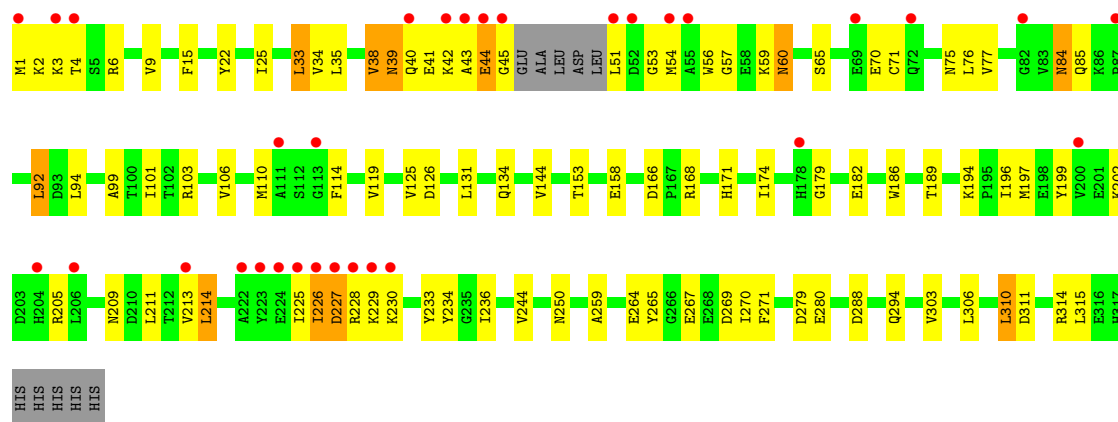


- Molecule 1: L-lactate dehydrogenase

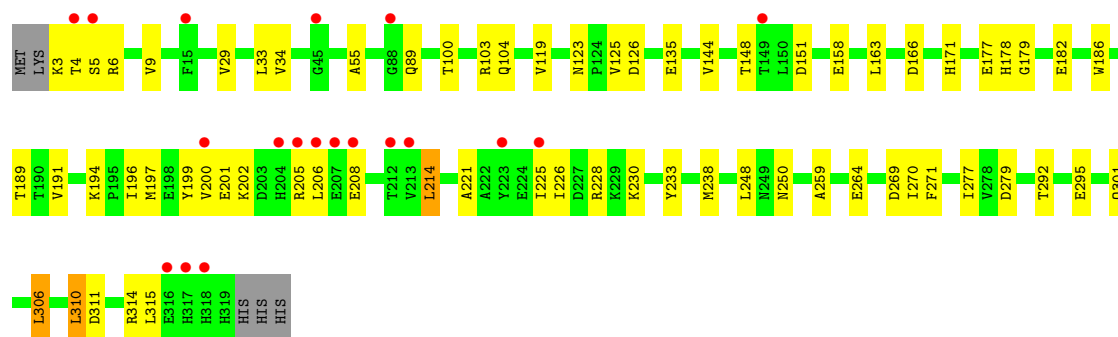
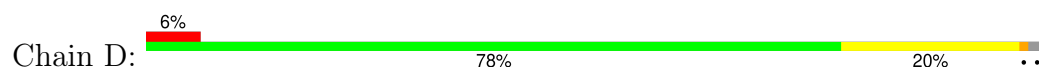


- Molecule 1: L-lactate dehydrogenase





● Molecule 1: L-lactate dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.43Å 127.82Å 133.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.96 – 2.38 35.94 – 2.38	Depositor EDS
% Data completeness (in resolution range)	96.4 (29.96-2.38) 99.5 (35.94-2.38)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.84 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.232 , 0.271 0.231 , 0.270	Depositor DCC
$R_{free}$ test set	3632 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.7	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 44.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.57$ , $\langle L^2 \rangle = 0.42$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10138	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.4886e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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	0.34	0/2436	0.58	0/3304
1	B	0.35	0/2480	0.59	0/3368
1	C	0.34	0/2436	0.56	1/3304 (0.0%)
1	D	0.34	0/2480	0.58	0/3368
All	All	0.34	0/9832	0.58	1/13344 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	174	ILE	N-CA-C	-5.13	97.16	111.00

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	2400	0	2415	83	0
1	B	2441	0	2442	55	0
1	C	2400	0	2415	76	0
1	D	2441	0	2442	54	0
2	A	12	0	16	2	0
2	B	12	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	18	0	24	0	0
2	D	6	0	8	0	0
3	A	86	0	0	1	0
3	B	121	0	0	3	0
3	C	79	0	0	1	0
3	D	122	0	0	3	0
All	All	10138	0	9778	253	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 13.

All (253) 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:250:ASN:HB2	1:A:280:GLU:HG3	1.57	0.87
1:B:226:ILE:HD11	1:B:230:LYS:HD2	1.55	0.86
1:C:158:GLU:HB3	1:C:214:LEU:HD11	1.57	0.85
1:A:2:LYS:HG3	1:C:75:ASN:ND2	1.96	0.81
1:C:38:VAL:HG23	1:C:39:ASN:H	1.47	0.80
1:A:6:ARG:HG2	1:C:4:THR:HB	1.65	0.78
1:A:38:VAL:O	1:A:39:ASN:HB2	1.84	0.77
1:A:217:LYS:O	1:A:221:ALA:HB2	1.83	0.77
1:B:92:LEU:HD22	1:B:96:LYS:HE3	1.68	0.75
1:A:85:GLN:HB2	1:A:94:LEU:HD22	1.67	0.75
1:A:200:VAL:HG23	1:A:201:GLU:HG3	1.69	0.74
1:D:103:ARG:HD2	1:D:315:LEU:HD21	1.69	0.74
1:A:39:ASN:HB3	1:A:41:GLU:HG2	1.70	0.72
1:D:177:GLU:OE1	1:D:306:LEU:HG	1.90	0.72
1:B:90:THR:O	1:B:91:ARG:HB2	1.91	0.70
1:C:166:ASP:OD2	1:C:168:ARG:HB2	1.92	0.69
1:B:200:VAL:HG22	1:B:206:LEU:HD12	1.73	0.69
1:D:89:GLN:CD	1:D:89:GLN:H	1.96	0.69
1:C:85:GLN:HB2	1:C:94:LEU:HD22	1.74	0.68
1:D:5:SER:HB2	3:D:592:HOH:O	1.94	0.67
1:D:226:ILE:HD11	1:D:230:LYS:HD2	1.77	0.67
1:C:310:LEU:HD22	1:C:314:ARG:HB3	1.76	0.67
1:C:34:VAL:HG11	1:C:70:GLU:HG3	1.78	0.66
1:C:158:GLU:HB3	1:C:214:LEU:CD1	2.25	0.66
1:C:51:LEU:HD21	1:C:65:SER:OG	1.96	0.66
1:A:250:ASN:OD1	1:A:279:ASP:HB2	1.95	0.65
1:A:76:LEU:HD13	1:A:244:VAL:HG13	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:ARG:O	1:D:206:LEU:HD12	1.96	0.65
1:A:4:THR:HB	1:C:6:ARG:HG2	1.79	0.65
1:C:42:LYS:HG3	1:C:43:ALA:H	1.61	0.64
1:B:221:ALA:HA	1:B:224:GLU:HG2	1.79	0.64
1:B:85:GLN:HE21	1:B:123:ASN:HB3	1.62	0.64
1:B:85:GLN:NE2	1:B:123:ASN:HB3	2.13	0.63
1:B:103:ARG:HD2	1:B:315:LEU:HD11	1.81	0.63
1:B:180:ASP:HA	1:B:219:LYS:HD2	1.80	0.63
1:D:151:ASP:OD1	1:D:178:HIS:HB2	1.98	0.63
1:C:250:ASN:OD1	1:C:280:GLU:HG2	1.99	0.62
1:A:3:LYS:HE3	3:C:667:HOH:O	1.99	0.62
1:B:15:PHE:CZ	1:B:44:GLU:HG2	2.33	0.62
1:C:134:GLN:HB3	1:C:315:LEU:HD12	1.80	0.62
1:B:200:VAL:HG22	1:B:206:LEU:CD1	2.30	0.61
1:C:103:ARG:NH2	1:C:131:LEU:HD22	2.16	0.61
1:B:5:SER:HB2	3:B:610:HOH:O	2.01	0.60
1:D:301:GLN:HG2	3:D:608:HOH:O	2.02	0.60
1:B:233:TYR:O	1:B:238:MET:HG3	2.01	0.60
1:D:103:ARG:HH11	1:D:103:ARG:HG2	1.67	0.60
1:B:151:ASP:OD1	1:B:178:HIS:HB2	2.02	0.59
1:D:186:TRP:HE3	1:D:196:ILE:HG21	1.67	0.58
1:A:250:ASN:CB	1:A:280:GLU:HG3	2.31	0.58
1:A:228:ARG:HG3	1:A:228:ARG:HH11	1.67	0.58
1:D:3:LYS:HG2	1:D:4:THR:H	1.68	0.58
1:D:29:VAL:HG12	1:D:248:LEU:HD12	1.86	0.58
1:B:197:MET:HA	1:B:197:MET:CE	2.34	0.58
1:D:100:THR:O	1:D:104:GLN:HG3	2.04	0.58
1:A:75:ASN:ND2	1:C:2:LYS:HB3	2.19	0.57
1:C:84:ASN:HB3	1:C:228:ARG:HH21	1.68	0.57
1:D:250:ASN:OD1	1:D:279:ASP:HB2	2.03	0.57
1:A:72:GLN:HG2	1:A:112:SER:O	2.05	0.57
1:A:39:ASN:CG	1:A:40:GLN:H	2.08	0.57
1:C:119:VAL:HA	1:C:144:VAL:O	2.05	0.56
1:C:186:TRP:HE3	1:C:196:ILE:HG21	1.70	0.56
1:D:179:GLY:O	1:D:182:GLU:HG2	2.06	0.55
1:A:267:GLU:HG3	1:A:303:VAL:HG21	1.88	0.55
1:A:215:ALA:O	1:A:219:LYS:HB2	2.05	0.55
1:A:223:TYR:HD1	1:A:223:TYR:O	1.89	0.55
1:C:269:ASP:O	1:C:270:ILE:HG23	2.07	0.55
1:B:67:THR:HG22	1:B:68:TYR:N	2.22	0.55
1:B:3:LYS:HB2	3:B:610:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:VAL:HG13	1:A:126:ASP:N	2.23	0.54
1:B:177:GLU:OE1	1:B:306:LEU:HG	2.08	0.54
1:A:228:ARG:HG2	1:A:233:TYR:OH	2.08	0.54
1:B:4:THR:HG21	1:D:248:LEU:HA	1.89	0.54
1:A:194:LYS:HD3	1:A:199:TYR:CE1	2.42	0.54
1:B:292:THR:OG1	1:B:295:GLU:HG3	2.08	0.54
1:B:100:THR:O	1:B:104:GLN:HG3	2.08	0.54
1:C:226:ILE:HG22	1:C:227:ASP:N	2.22	0.54
1:C:267:GLU:HG3	1:C:303:VAL:HG21	1.88	0.54
1:A:14:GLY:HA3	1:A:43:ALA:HB3	1.89	0.53
1:B:6:ARG:HG3	1:B:248:LEU:HD13	1.89	0.53
1:D:123:ASN:HA	1:D:125:VAL:N	2.24	0.53
1:A:223:TYR:O	1:A:223:TYR:CD1	2.62	0.53
1:C:250:ASN:OD1	1:C:279:ASP:HB2	2.08	0.53
1:A:228:ARG:HG3	1:A:228:ARG:NH1	2.23	0.53
1:A:260:TYR:CZ	1:A:269:ASP:HA	2.44	0.53
1:C:42:LYS:NZ	1:C:45:GLY:HA2	2.23	0.52
1:A:207:GLU:HG2	1:A:208:GLU:N	2.23	0.52
1:C:76:LEU:HD13	1:C:244:VAL:HG13	1.91	0.52
1:C:311:ASP:HA	1:C:314:ARG:HD2	1.90	0.52
1:A:198:GLU:HG2	1:A:202:LYS:HE2	1.92	0.52
1:C:270:ILE:HD12	1:C:270:ILE:O	2.10	0.52
1:D:264:GLU:H	1:D:264:GLU:CD	2.13	0.52
1:C:103:ARG:HH21	1:C:131:LEU:HD22	1.73	0.52
1:C:227:ASP:O	1:C:228:ARG:HB2	2.10	0.52
1:B:179:GLY:O	1:B:182:GLU:HG2	2.10	0.51
1:C:171:HIS:O	1:C:189:THR:HA	2.10	0.51
1:D:200:VAL:HG12	1:D:206:LEU:HB2	1.92	0.51
1:A:67:THR:OG1	1:A:69:GLU:HG2	2.10	0.51
2:A:402:GOL:H31	1:D:277:ILE:HD11	1.92	0.51
1:B:152:THR:HG22	2:B:401:GOL:H31	1.91	0.51
1:A:35:LEU:HD11	1:A:54:MET:HE1	1.93	0.51
1:A:221:ALA:C	1:A:223:TYR:H	2.14	0.51
1:A:264:GLU:HA	1:A:288:ASP:OD2	2.11	0.51
1:A:115:ASP:O	1:C:1:MET:HA	2.11	0.51
1:B:267:GLU:HG3	1:B:303:VAL:HG21	1.93	0.51
1:D:200:VAL:HG12	1:D:206:LEU:CB	2.41	0.51
1:B:9:VAL:HG22	1:B:34:VAL:HB	1.92	0.51
1:C:42:LYS:HG3	1:C:43:ALA:N	2.25	0.51
1:C:43:ALA:O	1:C:44:GLU:HB3	2.11	0.51
1:A:200:VAL:HG12	1:A:206:LEU:CB	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ASP:OD1	1:A:314:ARG:NH1	2.44	0.50
1:B:85:GLN:NE2	1:B:223:TYR:HE1	2.09	0.50
1:B:250:ASN:OD1	1:B:279:ASP:HB2	2.11	0.50
1:D:103:ARG:NH1	1:D:135:GLU:OE1	2.45	0.50
1:D:103:ARG:CD	1:D:315:LEU:HD21	2.40	0.50
1:A:158:GLU:HB3	1:A:214:LEU:CD2	2.42	0.50
1:A:179:GLY:O	1:A:182:GLU:HG2	2.11	0.50
1:C:179:GLY:O	1:C:182:GLU:HG2	2.11	0.50
1:D:270:ILE:O	1:D:270:ILE:HD12	2.12	0.49
1:A:34:VAL:HG21	1:A:70:GLU:HG2	1.93	0.49
1:B:125:VAL:HG13	1:B:126:ASP:N	2.27	0.49
1:B:310:LEU:HD22	1:B:314:ARG:HB3	1.93	0.49
1:A:209:ASN:O	1:A:213:VAL:HG23	2.13	0.49
1:C:42:LYS:CG	1:C:43:ALA:H	2.24	0.49
1:A:7:LYS:HZ1	1:A:70:GLU:HG3	1.78	0.49
1:B:95:VAL:HG22	1:B:124:PRO:HG3	1.93	0.49
1:A:186:TRP:HE3	1:A:196:ILE:HG21	1.79	0.48
1:C:84:ASN:CB	1:C:228:ARG:HH21	2.25	0.48
1:C:264:GLU:HA	1:C:288:ASP:OD2	2.14	0.48
1:D:3:LYS:HG2	1:D:4:THR:N	2.28	0.48
1:D:200:VAL:HG23	1:D:201:GLU:HG3	1.93	0.48
1:A:7:LYS:HB3	1:A:74:ALA:HA	1.95	0.48
1:B:194:LYS:HD3	1:B:199:TYR:CE1	2.49	0.48
1:A:15:PHE:HB3	1:A:233:TYR:CE2	2.49	0.48
1:D:148:THR:HA	3:D:517:HOH:O	2.13	0.48
1:A:42:LYS:HG3	1:A:44:GLU:H	1.78	0.48
1:C:15:PHE:CD2	1:C:43:ALA:HB1	2.49	0.48
1:B:206:LEU:HD12	1:B:206:LEU:O	2.14	0.48
1:C:228:ARG:HG3	1:C:228:ARG:HH11	1.79	0.48
1:C:41:GLU:HB3	1:C:101:ILE:HD13	1.96	0.47
1:C:230:LYS:HA	1:D:228:ARG:HG3	1.95	0.47
1:A:76:LEU:HD23	1:A:77:VAL:N	2.30	0.47
1:A:248:LEU:HD22	1:C:4:THR:HG21	1.97	0.47
1:C:33:LEU:HD13	1:C:35:LEU:HD21	1.95	0.47
1:A:54:MET:HE2	1:A:63:VAL:HG11	1.96	0.47
1:D:171:HIS:O	1:D:189:THR:HA	2.15	0.47
1:C:197:MET:HE2	1:C:211:LEU:HD11	1.95	0.47
1:C:233:TYR:O	1:C:236:ILE:HG22	2.15	0.47
1:D:194:LYS:HD3	1:D:199:TYR:CE1	2.50	0.47
1:D:225:ILE:HG23	1:D:225:ILE:O	2.14	0.47
1:D:310:LEU:HD22	1:D:314:ARG:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:LEU:HD23	1:A:76:LEU:C	2.35	0.47
1:B:264:GLU:CD	1:B:264:GLU:H	2.17	0.47
1:A:245:LYS:HE2	3:A:565:HOH:O	2.14	0.46
1:A:315:LEU:N	1:A:315:LEU:HD12	2.30	0.46
1:A:34:VAL:HG11	1:A:70:GLU:HG2	1.98	0.46
1:D:197:MET:O	1:D:200:VAL:HG22	2.15	0.46
1:B:160:ALA:HB1	1:B:165:VAL:O	2.16	0.46
1:B:121:ALA:HB1	1:B:236:ILE:HD11	1.98	0.46
1:D:311:ASP:HA	1:D:314:ARG:HD2	1.98	0.46
1:C:186:TRP:HB3	1:C:197:MET:HE3	1.98	0.46
1:C:71:CYS:O	1:C:114:PHE:HD1	1.99	0.46
1:C:59:LYS:O	1:C:60:ASN:C	2.54	0.46
1:A:233:TYR:O	1:A:236:ILE:HG22	2.17	0.45
1:D:205:ARG:C	1:D:206:LEU:HD12	2.37	0.45
1:B:213:VAL:O	1:B:217:LYS:HG3	2.17	0.45
1:D:233:TYR:O	1:D:238:MET:HG3	2.17	0.45
1:A:291:ILE:HB	1:A:295:GLU:OE1	2.16	0.45
1:C:153:THR:HG22	1:D:55:ALA:HB1	1.97	0.45
1:B:123:ASN:HA	1:B:125:VAL:N	2.32	0.44
1:B:63:VAL:O	1:B:64:TRP:HB3	2.16	0.44
1:C:158:GLU:CB	1:C:214:LEU:HD11	2.39	0.44
1:B:4:THR:HG21	1:D:248:LEU:HD23	1.98	0.44
1:C:9:VAL:HB	1:C:77:VAL:HG22	2.00	0.44
1:C:38:VAL:HG23	1:C:39:ASN:N	2.24	0.44
1:C:76:LEU:HD23	1:C:76:LEU:C	2.38	0.44
1:D:292:THR:OG1	1:D:295:GLU:HG3	2.18	0.44
1:B:270:ILE:HD12	1:B:270:ILE:O	2.17	0.44
1:C:39:ASN:HB3	1:C:40:GLN:H	1.51	0.44
1:A:251:GLU:O	1:A:252:GLN:HB2	2.18	0.44
1:B:186:TRP:HE3	1:B:196:ILE:HG21	1.82	0.44
1:C:3:LYS:HA	1:C:3:LYS:HD3	1.76	0.44
1:A:35:LEU:HD11	1:A:54:MET:CE	2.47	0.44
1:B:76:LEU:HD13	1:B:244:VAL:HG13	1.99	0.44
1:D:119:VAL:HA	1:D:144:VAL:O	2.17	0.44
1:C:76:LEU:HD23	1:C:77:VAL:N	2.32	0.43
1:A:149:THR:HA	1:A:256:PRO:HG2	1.99	0.43
1:C:25:ILE:CD1	1:C:54:MET:HG3	2.48	0.43
1:C:125:VAL:HG13	1:C:126:ASP:N	2.33	0.43
1:A:119:VAL:HA	1:A:144:VAL:O	2.18	0.43
1:C:44:GLU:OE2	1:C:229:LYS:HE2	2.17	0.43
1:A:171:HIS:O	1:A:189:THR:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:VAL:O	1:A:217:LYS:HG3	2.19	0.43
1:C:106:VAL:O	1:C:110:MET:HG2	2.18	0.43
1:A:11:VAL:HG11	1:A:105:ILE:CG2	2.49	0.43
1:D:125:VAL:HG13	1:D:126:ASP:N	2.34	0.43
1:A:15:PHE:N	1:A:43:ALA:HB3	2.34	0.43
1:A:259:ALA:O	1:A:271:PHE:HA	2.19	0.43
1:B:301:GLN:HG2	3:B:574:HOH:O	2.18	0.43
1:B:67:THR:CG2	1:B:68:TYR:N	2.82	0.43
1:C:33:LEU:CD1	1:C:35:LEU:HD21	2.49	0.43
1:C:259:ALA:O	1:C:271:PHE:HA	2.19	0.42
1:D:3:LYS:CG	1:D:4:THR:H	2.32	0.42
1:D:89:GLN:CD	1:D:89:GLN:N	2.67	0.42
1:C:264:GLU:O	1:C:265:TYR:HB2	2.20	0.42
1:A:145:GLY:HA3	1:A:258:SER:HB2	2.00	0.42
1:B:315:LEU:HD12	1:B:315:LEU:N	2.35	0.42
1:C:186:TRP:CE3	1:C:196:ILE:HG21	2.53	0.42
1:A:103:ARG:NE	1:A:315:LEU:HD11	2.35	0.42
1:A:200:VAL:HG12	1:A:206:LEU:HB3	2.01	0.42
1:C:99:ALA:HB1	1:C:131:LEU:CD1	2.49	0.42
1:D:196:ILE:O	1:D:199:TYR:HB2	2.19	0.42
2:A:402:GOL:H31	1:D:277:ILE:CD1	2.50	0.42
1:B:248:LEU:CD2	1:D:4:THR:HG21	2.50	0.42
1:A:240:THR:O	1:A:244:VAL:HG23	2.19	0.42
1:B:103:ARG:NH1	1:B:315:LEU:HG	2.35	0.42
1:D:269:ASP:O	1:D:270:ILE:HG23	2.20	0.42
1:A:4:THR:CG2	1:C:75:ASN:ND2	2.83	0.42
1:A:69:GLU:HG3	1:A:70:GLU:N	2.35	0.42
1:A:162:LYS:HA	1:A:162:LYS:HD2	1.87	0.42
1:A:226:ILE:HG22	1:A:227:ASP:N	2.33	0.42
1:C:42:LYS:CE	1:C:45:GLY:HA2	2.50	0.42
1:D:9:VAL:HG22	1:D:34:VAL:HB	2.00	0.42
1:A:216:ASP:HA	1:A:219:LYS:HB3	2.02	0.42
1:D:259:ALA:O	1:D:271:PHE:HA	2.19	0.42
1:C:54:MET:SD	1:C:54:MET:N	2.93	0.41
1:C:294:GLN:OE1	1:C:294:GLN:N	2.47	0.41
1:A:207:GLU:HG2	1:A:209:ASN:H	1.85	0.41
1:C:51:LEU:HD21	1:C:65:SER:CB	2.51	0.41
1:C:194:LYS:HD3	1:C:199:TYR:CE1	2.54	0.41
1:A:186:TRP:CE3	1:A:196:ILE:HG21	2.55	0.41
1:C:234:TYR:HE2	1:D:233:TYR:HH	1.66	0.41
1:A:142:ARG:HD3	1:C:1:MET:SD	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:MET:O	1:A:200:VAL:HG22	2.21	0.41
1:B:264:GLU:HG3	1:B:286:ILE:HG21	2.01	0.41
1:C:92:LEU:HD12	1:C:92:LEU:HA	1.91	0.41
1:C:209:ASN:O	1:C:213:VAL:HG23	2.20	0.41
1:A:39:ASN:ND2	1:A:40:GLN:H	2.19	0.41
1:A:236:ILE:HG23	1:A:237:GLY:N	2.35	0.41
1:A:264:GLU:O	1:A:265:TYR:HB2	2.21	0.41
1:A:315:LEU:N	1:A:315:LEU:CD1	2.84	0.41
1:D:158:GLU:HB2	1:D:214:LEU:HD11	2.03	0.41
1:A:9:VAL:HB	1:A:77:VAL:HG22	2.03	0.41
1:B:15:PHE:CD1	1:B:228:ARG:HD3	2.56	0.41
1:B:92:LEU:HD21	1:B:312:THR:OG1	2.21	0.40
1:B:145:GLY:HA3	1:B:258:SER:HB2	2.03	0.40
1:B:92:LEU:HD22	1:B:96:LYS:CE	2.46	0.40
1:B:92:LEU:O	1:B:96:LYS:HG3	2.22	0.40
1:B:149:THR:HA	1:B:256:PRO:HG2	2.03	0.40
1:D:163:LEU:HD12	1:D:191:VAL:HG21	2.03	0.40
1:C:22:TYR:CZ	1:D:230:LYS:HE2	2.57	0.40
1:D:208:GLU:OE1	1:D:208:GLU:HA	2.21	0.40
1:A:186:TRP:CZ3	1:A:211:LEU:HD12	2.57	0.40
1:A:194:LYS:HD3	1:A:199:TYR:CZ	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	308/322 (96%)	283 (92%)	19 (6%)	6 (2%)	8	9
1	B	315/322 (98%)	296 (94%)	16 (5%)	3 (1%)	15	21
1	C	308/322 (96%)	281 (91%)	17 (6%)	10 (3%)	4	3
1	D	315/322 (98%)	297 (94%)	16 (5%)	2 (1%)	25	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1246/1288 (97%)	1157 (93%)	68 (6%)	21 (2%)	9 10

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	60	ASN
1	C	39	ASN
1	B	90	THR
1	B	202	LYS
1	C	60	ASN
1	C	226	ILE
1	D	202	LYS
1	A	44	GLU
1	A	56	TRP
1	C	56	TRP
1	C	202	LYS
1	C	227	ASP
1	D	221	ALA
1	B	91	ARG
1	C	53	GLY
1	C	57	GLY
1	A	202	LYS
1	A	228	ARG
1	C	44	GLU
1	C	38	VAL

### 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	262/271 (97%)	258 (98%)	4 (2%)	65 79
1	B	266/271 (98%)	258 (97%)	8 (3%)	41 59
1	C	262/271 (97%)	254 (97%)	8 (3%)	40 57
1	D	266/271 (98%)	260 (98%)	6 (2%)	50 68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1056/1084 (97%)	1030 (98%)	26 (2%)	47 65

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	59	LYS
1	A	84	ASN
1	A	227	ASP
1	B	6	ARG
1	B	33	LEU
1	B	50	LEU
1	B	92	LEU
1	B	166	ASP
1	B	197	MET
1	B	306	LEU
1	B	310	LEU
1	C	33	LEU
1	C	84	ASN
1	C	92	LEU
1	C	205	ARG
1	C	214	LEU
1	C	225	ILE
1	C	306	LEU
1	C	310	LEU
1	D	6	ARG
1	D	33	LEU
1	D	166	ASP
1	D	214	LEU
1	D	306	LEU
1	D	310	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	252	GLN
1	B	27	GLN
1	B	72	GLN
1	B	85	GLN
1	B	220	ASN

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Mol	Chain	Res	Type
1	B	318	HIS
1	C	85	GLN
1	D	72	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](#)

8 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	402	-	5,5,5	0.42	0	5,5,5	0.27	0
2	GOL	A	401	-	5,5,5	0.33	0	5,5,5	0.29	0
2	GOL	C	503	-	5,5,5	0.45	0	5,5,5	0.28	0
2	GOL	D	401	-	5,5,5	0.40	0	5,5,5	0.32	0
2	GOL	A	402	-	5,5,5	0.37	0	5,5,5	0.24	0
2	GOL	C	501	-	5,5,5	0.38	0	5,5,5	0.25	0
2	GOL	B	401	-	5,5,5	0.45	0	5,5,5	0.24	0
2	GOL	C	502	-	5,5,5	0.34	0	5,5,5	0.29	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	GOL	B	402	-	-	2/4/4/4	-
2	GOL	A	401	-	-	4/4/4/4	-
2	GOL	C	503	-	-	2/4/4/4	-
2	GOL	D	401	-	-	4/4/4/4	-
2	GOL	A	402	-	-	0/4/4/4	-
2	GOL	C	501	-	-	2/4/4/4	-
2	GOL	B	401	-	-	2/4/4/4	-
2	GOL	C	502	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	GOL	O1-C1-C2-O2
2	A	401	GOL	O1-C1-C2-C3
2	A	401	GOL	C1-C2-C3-O3
2	B	402	GOL	C1-C2-C3-O3
2	B	402	GOL	O2-C2-C3-O3
2	C	501	GOL	O1-C1-C2-C3
2	D	401	GOL	C1-C2-C3-O3
2	A	401	GOL	O2-C2-C3-O3
2	B	401	GOL	O1-C1-C2-C3
2	C	502	GOL	O1-C1-C2-C3
2	C	502	GOL	C1-C2-C3-O3
2	C	503	GOL	O1-C1-C2-C3
2	D	401	GOL	O1-C1-C2-C3
2	C	501	GOL	O1-C1-C2-O2
2	D	401	GOL	O2-C2-C3-O3
2	B	401	GOL	O1-C1-C2-O2
2	C	502	GOL	O1-C1-C2-O2
2	C	502	GOL	O2-C2-C3-O3
2	C	503	GOL	O1-C1-C2-O2
2	D	401	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	402	GOL	2	0
2	B	401	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

### 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, 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	312/322 (96%)	0.70	35 (11%) <b>5</b> <b>6</b>	33, 53, 108, 121	0
1	B	317/322 (98%)	0.33	17 (5%) <b>25</b> <b>28</b>	32, 49, 82, 98	0
1	C	312/322 (96%)	0.64	32 (10%) <b>6</b> <b>7</b>	33, 53, 110, 132	0
1	D	317/322 (98%)	0.29	19 (5%) <b>21</b> <b>24</b>	32, 49, 86, 99	0
All	All	1258/1288 (97%)	0.49	103 (8%) <b>11</b> <b>12</b>	32, 51, 96, 132	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	225	ILE	10.4
1	A	54	MET	7.2
1	C	43	ALA	7.0
1	A	225	ILE	6.7
1	C	45	GLY	6.7
1	A	227	ASP	6.6
1	A	55	ALA	6.3
1	C	223	TYR	6.2
1	C	44	GLU	6.0
1	A	317	HIS	5.9
1	D	318	HIS	5.9
1	B	225	ILE	5.8
1	C	222	ALA	5.8
1	C	4	THR	5.5
1	A	223	TYR	5.5
1	A	226	ILE	5.4
1	C	55	ALA	5.1
1	C	54	MET	4.9
1	A	4	THR	4.8
1	A	224	GLU	4.8
1	D	200	VAL	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	200	VAL	4.4
1	A	87	PRO	4.4
1	A	222	ALA	4.3
1	C	42	LYS	4.2
1	D	204	HIS	4.2
1	A	316	GLU	4.1
1	D	317	HIS	4.1
1	C	204	HIS	4.1
1	C	224	GLU	4.0
1	A	42	LYS	4.0
1	C	226	ILE	3.9
1	A	229	LYS	3.8
1	A	228	ARG	3.8
1	A	44	GLU	3.8
1	C	228	ARG	3.7
1	D	225	ILE	3.7
1	B	223	TYR	3.7
1	C	40	GLN	3.7
1	B	43	ALA	3.6
1	A	86	LYS	3.6
1	A	51	LEU	3.5
1	C	51	LEU	3.5
1	C	200	VAL	3.4
1	A	111	ALA	3.4
1	D	15	PHE	3.4
1	A	43	ALA	3.3
1	C	227	ASP	3.3
1	B	319	HIS	3.3
1	B	15	PHE	3.2
1	C	87	PRO	3.2
1	B	208	GLU	3.2
1	A	230	LYS	3.2
1	C	3	LYS	3.1
1	A	40	GLN	3.1
1	A	204	HIS	3.1
1	D	208	GLU	3.1
1	C	229	LYS	3.1
1	D	4	THR	3.0
1	B	204	HIS	3.0
1	A	52	ASP	2.9
1	D	223	TYR	2.9
1	A	69	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	206	LEU	2.9
1	B	4	THR	2.8
1	C	230	LYS	2.8
1	A	220	ASN	2.7
1	A	88	GLY	2.7
1	D	207	GLU	2.7
1	B	201	GLU	2.6
1	C	111	ALA	2.6
1	A	178	HIS	2.6
1	A	72	GLN	2.5
1	C	1	MET	2.5
1	D	316	GLU	2.4
1	C	72	GLN	2.4
1	C	69	GLU	2.4
1	D	213	VAL	2.4
1	D	5	SER	2.4
1	C	206	LEU	2.4
1	A	1	MET	2.4
1	B	207	GLU	2.4
1	C	52	ASP	2.4
1	C	178	HIS	2.3
1	A	53	GLY	2.3
1	C	113	GLY	2.3
1	C	82	GLY	2.3
1	B	42	LYS	2.3
1	D	88	GLY	2.3
1	B	226	ILE	2.2
1	D	149	THR	2.2
1	D	212	THR	2.2
1	B	209	ASN	2.2
1	B	148	THR	2.2
1	A	315	LEU	2.2
1	B	200	VAL	2.1
1	C	213	VAL	2.1
1	A	206	LEU	2.1
1	B	88	GLY	2.1
1	D	45	GLY	2.1
1	B	318	HIS	2.0
1	D	205	ARG	2.0
1	A	64	TRP	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 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( $\text{\AA}^2$ )	Q<0.9
2	GOL	C	502	6/6	0.60	0.24	95,98,98,99	0
2	GOL	C	503	6/6	0.64	0.48	80,81,82,82	0
2	GOL	B	402	6/6	0.71	0.32	86,87,88,88	0
2	GOL	C	501	6/6	0.71	0.37	90,92,92,92	0
2	GOL	A	401	6/6	0.76	0.32	91,93,93,94	0
2	GOL	B	401	6/6	0.83	0.32	57,62,63,64	0
2	GOL	A	402	6/6	0.90	0.16	74,77,77,77	0
2	GOL	D	401	6/6	0.90	0.24	56,60,62,62	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.