



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:18 PM GMT

PDB ID : 1JPM  
Title : L-Ala-D/L-Glu Epimerase  
Authors : Gulick, A.M.; Schmidt, D.M.Z.; Gerlt, J.A.; Rayment, I.  
Deposited on : 2001-08-02  
Resolution : 2.25 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  
<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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

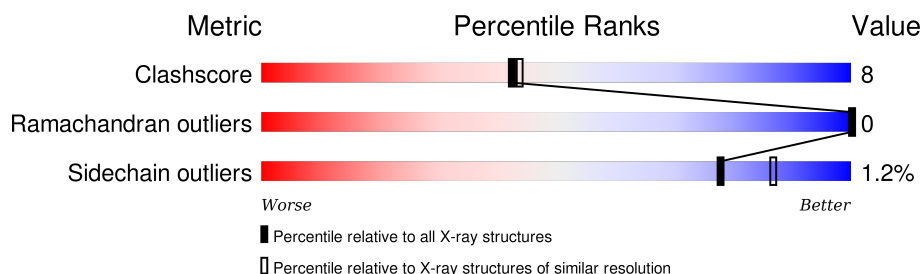
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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))
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	366	 83% 15% •
1	B	366	 81% 16% ••
1	C	366	 80% 18% •
1	D	366	 84% 13% •

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11271 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-Ala-D/L-Glu Epimerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	0	0
			2689	1697	459	515	18			
1	B	359	Total	C	N	O	S	0	0	0
			2685	1693	457	517	18			
1	C	358	Total	C	N	O	S	0	0	0
			2686	1695	457	516	18			
1	D	358	Total	C	N	O	S	0	0	0
			2681	1693	457	513	18			

- 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<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 4 is water.

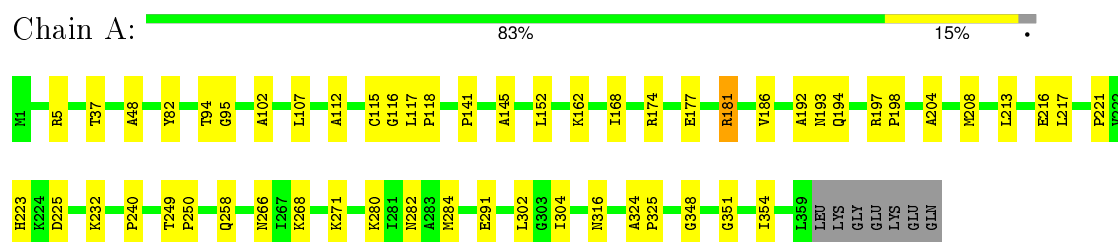
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	133	Total	O	0	0
			133	133		
4	B	157	Total	O	0	0
			157	157		
4	C	115	Total	O	0	0
			115	115		
4	D	115	Total	O	0	0
			115	115		

### 3 Residue-property plots

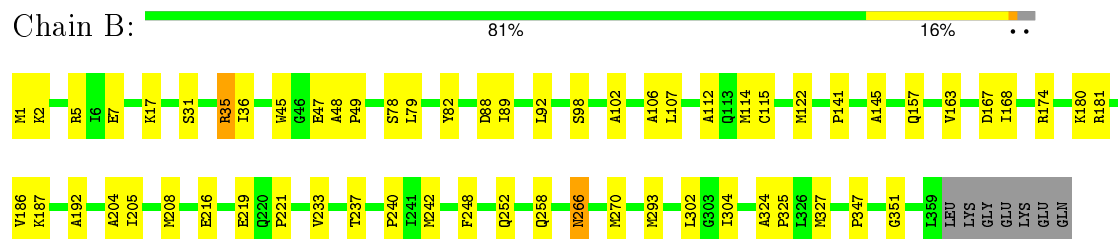
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

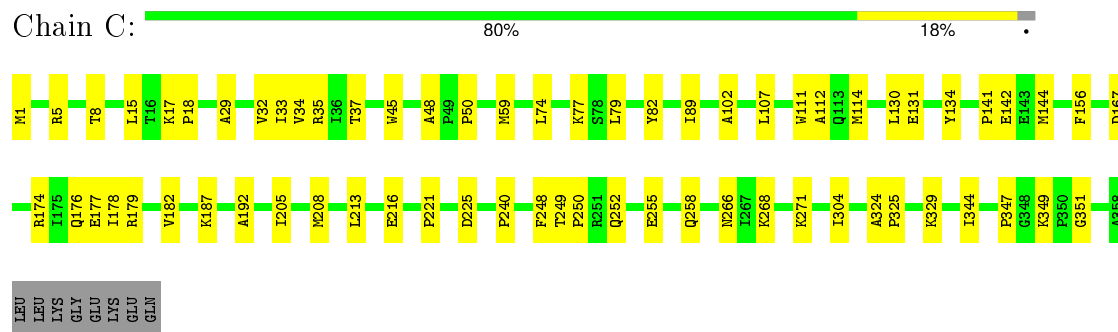
#### • Molecule 1: L-Ala-D/L-Glu Epimerase



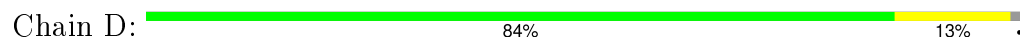
#### • Molecule 1: L-Ala-D/L-Glu Epimerase



#### • Molecule 1: L-Ala-D/L-Glu Epimerase



#### • Molecule 1: L-Ala-D/L-Glu Epimerase



R1	R2	R3	S31	V32	D39	W45	A48	P49	P50	M59	L74	I85	I89	M97	A102	L107	A112	A145	M150	K153	I168	Q176	E177	R181	R189	A192	I205	M208	L213	E216	P221	P240					
P248	Q252	N266	G273	N282	I298	L302	G303	I304	N316	A324	P325	T330	D331	N334	P347	G351	I354	I355	A358	LEU	LEU	LYS	GLY	GLU	LYS	GLU	GLN										

## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.63Å 157.63Å 168.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.25	Depositor
% Data completeness (in resolution range)	99.9 (30.00-2.25)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.197 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11271	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

## 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, 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	0.32	0/2731	0.58	0/3696
1	B	0.32	0/2727	0.58	0/3693
1	C	0.32	0/2728	0.58	0/3693
1	D	0.31	0/2723	0.58	0/3686
All	All	0.32	0/10909	0.58	0/14768

There are no bond length outliers.

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	2689	0	2725	40	0
1	B	2685	0	2713	53	0
1	C	2686	0	2729	49	0
1	D	2681	0	2722	42	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	6	0	8	0	0

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	133	0	0	2	0
4	B	157	0	0	3	0
4	C	115	0	0	4	0
4	D	115	0	0	2	0
All	All	11271	0	10897	180	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 8.

All (180) 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:49:PRO:HD2	1:B:270:MET:HE2	1.31	1.09
1:D:32:VAL:HG23	1:D:59:MET:HE2	1.50	0.94
1:A:162:LYS:HD3	1:A:193:ASN:HD22	1.41	0.85
1:D:168:ILE:HD12	1:D:168:ILE:H	1.42	0.83
1:C:248:PHE:H	1:C:252:GLN:NE2	1.79	0.81
1:B:115:CYS:HB2	1:B:122:MET:HE1	1.62	0.80
1:C:32:VAL:HG23	1:C:59:MET:HE2	1.67	0.77
1:C:347:PRO:HG2	4:C:1019:HOH:O	1.85	0.76
1:B:48:ALA:HB2	1:B:102:ALA:HB2	1.67	0.76
1:B:49:PRO:CD	1:B:270:MET:HE2	2.16	0.74
1:D:302:LEU:HD13	1:D:354:ILE:HD13	1.67	0.74
1:D:347:PRO:HG2	4:D:1023:HOH:O	1.88	0.74
1:B:5:ARG:HD3	1:B:7:GLU:OE1	1.87	0.74
1:D:282:ASN:HD21	1:D:316:ASN:HB3	1.52	0.73
1:C:248:PHE:H	1:C:252:GLN:HE21	1.38	0.71
1:D:45:TRP:CG	1:D:354:ILE:HD11	2.26	0.70
1:A:162:LYS:CD	1:A:193:ASN:HD22	2.03	0.70
1:A:162:LYS:HD3	1:A:193:ASN:ND2	2.06	0.70
1:D:150:ASN:O	1:D:153:LYS:HG2	1.91	0.69
1:B:233:VAL:O	1:B:237:THR:HG22	1.93	0.69
1:C:50:PRO:HD3	1:C:59:MET:CE	2.23	0.68
1:A:162:LYS:HE2	1:A:193:ASN:HB2	1.73	0.68
1:C:304:ILE:HG13	1:C:325:PRO:CG	2.25	0.67
1:B:180:LYS:HB2	1:B:180:LYS:NZ	2.11	0.66
1:C:48:ALA:HB2	1:C:102:ALA:HB2	1.79	0.64
1:C:255:GLU:HA	1:C:258:GLN:HE21	1.62	0.64
1:A:48:ALA:HB2	1:A:102:ALA:HB2	1.80	0.64
1:B:115:CYS:HB2	1:B:122:MET:CE	2.29	0.63
1:A:282:ASN:HD21	1:A:316:ASN:HB3	1.63	0.63

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:168:ILE:HD12	1:D:168:ILE:N	2.14	0.62
1:B:115:CYS:CB	1:B:122:MET:HE1	2.30	0.62
1:B:112:ALA:HB3	1:B:351:GLY:HA2	1.80	0.62
1:B:168:ILE:HD13	1:B:204:ALA:HB2	1.82	0.62
1:D:74:LEU:HD22	1:D:85:ILE:HG23	1.82	0.62
1:D:168:ILE:CD1	1:D:168:ILE:H	2.13	0.61
1:B:145:ALA:HB1	1:B:181:ARG:HG3	1.82	0.61
1:D:145:ALA:HB1	1:D:181:ARG:HG3	1.83	0.61
1:C:144:MET:SD	1:C:174:ARG:HD2	2.40	0.61
1:D:31:SER:HA	1:D:59:MET:HE1	1.83	0.61
1:C:268:LYS:HB2	1:C:271:LYS:HE2	1.82	0.60
1:B:1:MET:HG2	1:B:79:LEU:HD12	1.84	0.60
1:D:145:ALA:HB1	1:D:181:ARG:CG	2.32	0.60
1:A:304:ILE:HG13	1:A:325:PRO:HD3	1.84	0.60
1:D:304:ILE:HG13	1:D:325:PRO:CG	2.31	0.59
1:B:141:PRO:HB3	1:B:174:ARG:HA	1.84	0.58
1:A:162:LYS:CE	1:A:193:ASN:HD22	2.16	0.58
1:C:111:TRP:HD1	1:C:114:MET:HE2	1.68	0.58
1:A:118:PRO:HG3	1:A:348:GLY:O	2.04	0.57
1:B:180:LYS:HB2	1:B:180:LYS:HZ2	1.67	0.57
1:A:141:PRO:HB3	1:A:174:ARG:HA	1.85	0.57
1:D:355:ILE:N	1:D:355:ILE:HD12	2.20	0.57
1:C:248:PHE:HB2	1:C:252:GLN:HE22	1.68	0.57
1:C:8:THR:HG22	1:C:34:VAL:HG22	1.85	0.57
1:D:48:ALA:HB2	1:D:102:ALA:HB2	1.87	0.55
1:A:324:ALA:HB3	1:A:325:PRO:HD3	1.88	0.55
1:C:32:VAL:HG23	1:C:59:MET:CE	2.35	0.55
1:B:115:CYS:SG	1:B:122:MET:CE	2.95	0.55
1:B:327:MET:HE1	4:B:1036:HOH:O	2.07	0.55
1:A:268:LYS:HB2	1:A:271:LYS:HE2	1.89	0.55
1:D:304:ILE:HG13	1:D:325:PRO:HD3	1.88	0.55
1:B:347:PRO:HG2	4:B:1032:HOH:O	2.06	0.55
1:D:324:ALA:HB3	1:D:325:PRO:HD3	1.88	0.54
1:B:115:CYS:SG	1:B:122:MET:HE1	2.47	0.54
1:C:192:ALA:HB3	1:C:221:PRO:HA	1.90	0.54
1:B:145:ALA:HB1	1:B:181:ARG:CG	2.37	0.54
1:C:141:PRO:HB3	1:C:174:ARG:HA	1.89	0.53
1:D:89:ILE:HD11	1:D:107:LEU:CD1	2.39	0.53
1:D:50:PRO:HD3	1:D:59:MET:CE	2.39	0.53
1:A:145:ALA:HB1	1:A:181:ARG:HG3	1.90	0.53
1:B:48:ALA:HB1	1:B:270:MET:HE1	1.91	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:205:ILE:HA	1:C:208:MET:HE2	1.91	0.53
1:B:49:PRO:HD2	1:B:270:MET:CE	2.22	0.52
1:C:50:PRO:HD3	1:C:59:MET:HE1	1.89	0.52
1:A:192:ALA:HB3	1:A:221:PRO:HA	1.90	0.52
1:A:116:GLY:HA3	1:C:37:THR:HG21	1.92	0.52
1:A:112:ALA:HB3	1:A:351:GLY:HA2	1.92	0.52
1:A:304:ILE:HG13	1:A:325:PRO:CG	2.40	0.52
1:C:304:ILE:HG13	1:C:325:PRO:CD	2.41	0.51
1:D:298:ILE:HG12	4:D:1067:HOH:O	2.10	0.51
1:A:152:LEU:HD13	1:A:186:VAL:CG1	2.41	0.51
1:B:187:LYS:HA	4:B:1048:HOH:O	2.11	0.51
1:C:15:LEU:HD21	1:C:29:ALA:HB2	1.93	0.50
1:B:157:GLN:O	1:B:186:VAL:HA	2.11	0.50
1:C:35:ARG:HD3	1:C:45:TRP:CH2	2.46	0.50
1:A:304:ILE:HG13	1:A:325:PRO:CD	2.41	0.50
1:A:198:PRO:HB2	1:A:232:LYS:HD3	1.94	0.50
1:A:208:MET:HG2	1:A:213:LEU:HD12	1.94	0.50
1:D:304:ILE:HG13	1:D:325:PRO:CD	2.42	0.50
1:D:216:GLU:O	1:D:240:PRO:HG2	2.12	0.49
1:D:176:GLN:HG3	1:D:213:LEU:HD21	1.95	0.49
1:B:242:MET:HE3	1:B:293:MET:HE2	1.94	0.49
1:D:248:PHE:HB2	1:D:252:GLN:OE1	2.13	0.49
1:C:111:TRP:CD1	1:C:114:MET:HE2	2.48	0.48
1:B:304:ILE:HG13	1:B:325:PRO:HD3	1.96	0.48
1:B:304:ILE:HG13	1:B:325:PRO:CG	2.44	0.48
1:A:280:LYS:O	1:A:284:MET:HG3	2.13	0.48
1:C:304:ILE:HG13	1:C:325:PRO:HD3	1.96	0.48
1:B:324:ALA:HB3	1:B:325:PRO:HD3	1.95	0.47
1:D:192:ALA:HB3	1:D:221:PRO:HA	1.96	0.47
1:B:112:ALA:HA	1:B:122:MET:SD	2.55	0.47
1:D:145:ALA:HB2	1:D:177:GLU:HG3	1.95	0.47
1:B:242:MET:CE	1:B:266:ASN:HB2	2.44	0.47
1:A:197:ARG:NH1	4:A:2044:HOH:O	2.48	0.47
1:C:59:MET:HA	1:C:59:MET:CE	2.45	0.47
1:C:205:ILE:HA	1:C:208:MET:CE	2.44	0.47
1:A:115:CYS:O	1:C:5:ARG:NH2	2.47	0.47
1:B:192:ALA:HB3	1:B:221:PRO:HA	1.97	0.47
1:B:115:CYS:CB	1:B:122:MET:CE	2.91	0.46
1:B:98:SER:HB3	1:B:270:MET:CE	2.45	0.46
1:B:88:ASP:O	1:B:92:LEU:HG	2.15	0.46
1:D:3:ILE:HD11	1:D:107:LEU:HD21	1.96	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:PRO:HD3	1:C:59:MET:HE3	1.96	0.46
1:C:112:ALA:HB3	1:C:351:GLY:HA2	1.98	0.46
1:B:115:CYS:SG	1:B:122:MET:HE2	2.56	0.46
1:D:45:TRP:CD2	1:D:354:ILE:HD11	2.50	0.46
1:A:145:ALA:HB2	1:A:177:GLU:HG3	1.98	0.45
1:B:205:ILE:HA	1:B:208:MET:CE	2.46	0.45
1:C:1:MET:CE	1:C:79:LEU:HD22	2.46	0.45
1:D:330:THR:HG22	1:D:331:ASP:N	2.30	0.45
1:C:176:GLN:HA	1:C:213:LEU:HD22	1.97	0.45
1:A:181:ARG:CG	1:A:181:ARG:HH11	2.29	0.45
1:D:205:ILE:HA	1:D:208:MET:CE	2.47	0.45
1:D:355:ILE:H	1:D:355:ILE:HD12	1.79	0.45
1:A:217:LEU:HD12	1:A:217:LEU:C	2.36	0.45
1:A:168:ILE:HD13	1:A:204:ALA:HB2	1.98	0.45
1:D:112:ALA:HB3	1:D:351:GLY:HA2	1.98	0.45
1:B:219:GLU:HG2	1:B:242:MET:CE	2.47	0.45
1:A:216:GLU:O	1:A:240:PRO:HG2	2.16	0.45
1:C:35:ARG:HD3	1:C:45:TRP:CZ3	2.52	0.44
1:C:349:LYS:HG3	4:C:1074:HOH:O	2.17	0.44
1:C:59:MET:HE2	1:C:59:MET:HA	2.00	0.44
1:C:8:THR:HA	1:C:33:ILE:O	2.18	0.44
1:C:216:GLU:O	1:C:240:PRO:HG2	2.18	0.44
1:A:82:TYR:CE1	1:A:107:LEU:HG	2.53	0.44
1:D:145:ALA:HB1	1:D:181:ARG:HG2	2.00	0.43
1:B:98:SER:HB3	1:B:270:MET:HE1	2.00	0.43
1:A:302:LEU:HD13	1:A:354:ILE:HG13	2.01	0.43
1:C:134:TYR:HB2	1:C:156:PHE:CD2	2.54	0.43
1:C:141:PRO:O	1:C:177:GLU:HG2	2.18	0.43
1:A:94:THR:HG22	1:A:95:GLY:N	2.34	0.43
1:D:2:LYS:HB2	1:D:39:ASP:OD2	2.18	0.43
1:B:181:ARG:HH11	1:B:181:ARG:HG2	1.84	0.43
1:D:205:ILE:HA	1:D:208:MET:HE3	2.01	0.43
1:B:35:ARG:HD2	1:B:45:TRP:CZ3	2.54	0.43
1:D:354:ILE:O	1:D:354:ILE:HG13	2.19	0.42
1:D:334:ASN:HD22	1:D:355:ILE:HB	1.84	0.42
1:C:74:LEU:O	1:C:77:LYS:HB2	2.19	0.42
1:B:31:SER:HA	1:B:49:PRO:HA	2.00	0.42
1:C:15:LEU:O	1:C:329:LYS:HE3	2.19	0.42
1:C:179:ARG:NE	1:C:213:LEU:O	2.50	0.42
1:B:216:GLU:O	1:B:240:PRO:HG2	2.19	0.42
1:B:248:PHE:HB2	1:B:252:GLN:OE1	2.19	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:178:ILE:O	1:C:182:VAL:HG22	2.19	0.42
1:C:89:ILE:HD11	1:C:107:LEU:HD11	2.01	0.42
1:C:324:ALA:HB3	1:C:325:PRO:HD3	2.02	0.42
1:C:131:GLU:HG3	4:C:1118:HOH:O	2.19	0.42
1:B:304:ILE:HG13	1:B:325:PRO:CD	2.49	0.42
1:D:97:MET:O	1:D:273:GLY:HA2	2.19	0.42
1:A:258:GLN:HG3	1:B:258:GLN:HG3	2.02	0.42
1:C:187:LYS:HA	4:C:1034:HOH:O	2.20	0.42
1:D:282:ASN:ND2	1:D:316:ASN:HD22	2.18	0.42
1:B:89:ILE:HD11	1:B:107:LEU:CD1	2.49	0.42
1:C:130:LEU:CD1	1:C:344:ILE:HB	2.50	0.42
1:C:17:LYS:HA	1:C:18:PRO:HD2	1.81	0.42
1:B:2:LYS:HG2	1:B:78:SER:HB3	2.01	0.41
1:D:189:ARG:HG3	1:D:216:GLU:OE2	2.20	0.41
1:B:47:GLU:OE1	1:B:302:LEU:HB3	2.20	0.41
1:A:223:HIS:HD2	1:A:225:ASP:H	1.69	0.41
1:A:5:ARG:HB3	1:A:37:THR:HB	2.03	0.41
1:A:152:LEU:HD13	1:A:186:VAL:HG12	2.02	0.41
1:B:89:ILE:HD11	1:B:107:LEU:HD11	2.02	0.41
1:A:249:THR:HB	1:A:250:PRO:HD2	2.03	0.41
1:B:36:ILE:HD12	1:B:106:ALA:HB3	2.03	0.41
1:A:194:GLN:HB3	1:A:223:HIS:HA	2.03	0.41
1:B:163:VAL:O	1:B:192:ALA:HA	2.21	0.41
1:A:291:GLU:HG3	4:A:2106:HOH:O	2.21	0.41
1:D:304:ILE:HG13	1:D:325:PRO:HG3	2.02	0.40
1:B:242:MET:HE3	1:B:293:MET:CE	2.51	0.40
1:C:249:THR:HB	1:C:250:PRO:HD2	2.03	0.40
1:A:117:LEU:HD11	1:B:114:MET:HE3	2.04	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	357/366 (98%)	343 (96%)	14 (4%)	0	100	100
1	B	357/366 (98%)	339 (95%)	18 (5%)	0	100	100
1	C	356/366 (97%)	341 (96%)	15 (4%)	0	100	100
1	D	356/366 (97%)	341 (96%)	15 (4%)	0	100	100
All	All	1426/1464 (97%)	1364 (96%)	62 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	280/292 (96%)	278 (99%)	2 (1%)	88	93
1	B	280/292 (96%)	275 (98%)	5 (2%)	66	77
1	C	282/292 (97%)	277 (98%)	5 (2%)	66	77
1	D	280/292 (96%)	279 (100%)	1 (0%)	93	96
All	All	1122/1168 (96%)	1109 (99%)	13 (1%)	78	87

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	181	ARG
1	A	266	ASN
1	B	17	LYS
1	B	35	ARG
1	B	82	TYR
1	B	167	ASP
1	B	266	ASN
1	C	82	TYR
1	C	142	GLU
1	C	167	ASP
1	C	225	ASP
1	C	266	ASN
1	D	266	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	176	GLN
1	A	193	ASN
1	A	223	HIS
1	A	258	GLN
1	A	266	ASN
1	A	282	ASN
1	A	334	ASN
1	B	68	HIS
1	B	176	GLN
1	B	258	GLN
1	B	266	ASN
1	B	334	ASN
1	C	90	GLN
1	C	91	HIS
1	C	150	ASN
1	C	157	GLN
1	C	176	GLN
1	C	252	GLN
1	C	258	GLN
1	C	266	ASN
1	D	176	GLN
1	D	258	GLN
1	D	266	ASN
1	D	282	ASN
1	D	334	ASN

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 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	2001	-	5,5,5	0.92	0	5,5,5	0.38	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
3	GOL	A	2001	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.