



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2024 – 10:34 PM EST

PDB ID : 8GQL  
Title : X-ray structure of Thiolase from Pseudomonas aeruginosa PAO1  
Authors : Hong, J.; Son, H.F.; Kim, K.J.  
Deposited on : 2022-08-30  
Resolution : 2.30 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 : 2.36  
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

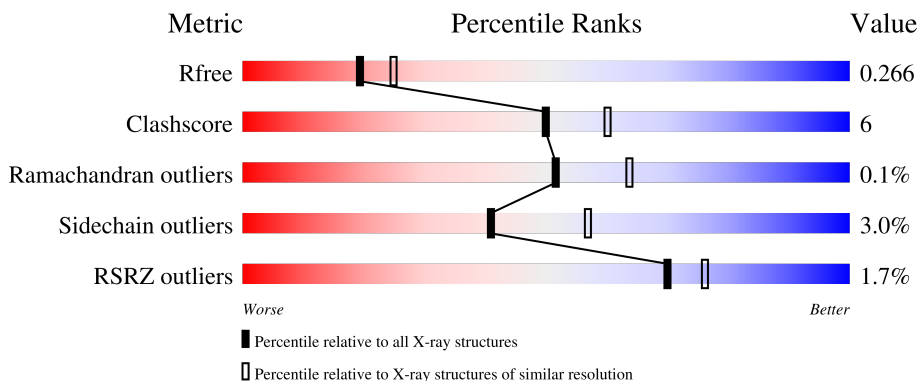
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	399	
1	B	399	
1	C	399	
1	D	399	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 11523 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 Thiolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	390	2822	1747	508	546	21	0	0	0
1	B	387	2796	1732	501	542	21	0	0	0
1	C	386	2789	1728	500	540	21	0	0	0
1	D	386	2789	1728	500	540	21	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

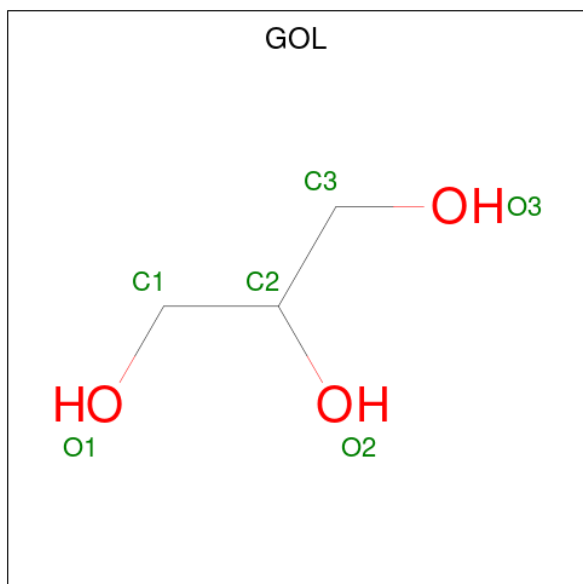
Chain	Residue	Modelled	Actual	Comment	Reference
A	392	LEU	-	expression tag	UNP G3XD40
A	393	GLU	-	expression tag	UNP G3XD40
A	394	HIS	-	expression tag	UNP G3XD40
A	395	HIS	-	expression tag	UNP G3XD40
A	396	HIS	-	expression tag	UNP G3XD40
A	397	HIS	-	expression tag	UNP G3XD40
A	398	HIS	-	expression tag	UNP G3XD40
A	399	HIS	-	expression tag	UNP G3XD40
B	392	LEU	-	expression tag	UNP G3XD40
B	393	GLU	-	expression tag	UNP G3XD40
B	394	HIS	-	expression tag	UNP G3XD40
B	395	HIS	-	expression tag	UNP G3XD40
B	396	HIS	-	expression tag	UNP G3XD40
B	397	HIS	-	expression tag	UNP G3XD40
B	398	HIS	-	expression tag	UNP G3XD40
B	399	HIS	-	expression tag	UNP G3XD40
C	392	LEU	-	expression tag	UNP G3XD40
C	393	GLU	-	expression tag	UNP G3XD40
C	394	HIS	-	expression tag	UNP G3XD40
C	395	HIS	-	expression tag	UNP G3XD40
C	396	HIS	-	expression tag	UNP G3XD40

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Chain	Residue	Modelled	Actual	Comment	Reference
C	397	HIS	-	expression tag	UNP G3XD40
C	398	HIS	-	expression tag	UNP G3XD40
C	399	HIS	-	expression tag	UNP G3XD40
D	392	LEU	-	expression tag	UNP G3XD40
D	393	GLU	-	expression tag	UNP G3XD40
D	394	HIS	-	expression tag	UNP G3XD40
D	395	HIS	-	expression tag	UNP G3XD40
D	396	HIS	-	expression tag	UNP G3XD40
D	397	HIS	-	expression tag	UNP G3XD40
D	398	HIS	-	expression tag	UNP G3XD40
D	399	HIS	-	expression tag	UNP G3XD40

- 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	C	1	Total	C O	0	0
			6	3 3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	84	Total	O	0	0
			84	84		
3	C	78	Total	O	0	0
			78	78		

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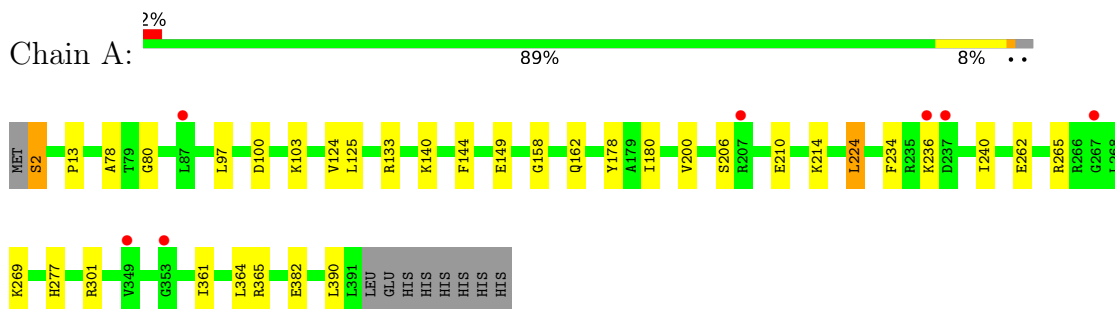
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	D	59	Total	O	0	0
			59	59		

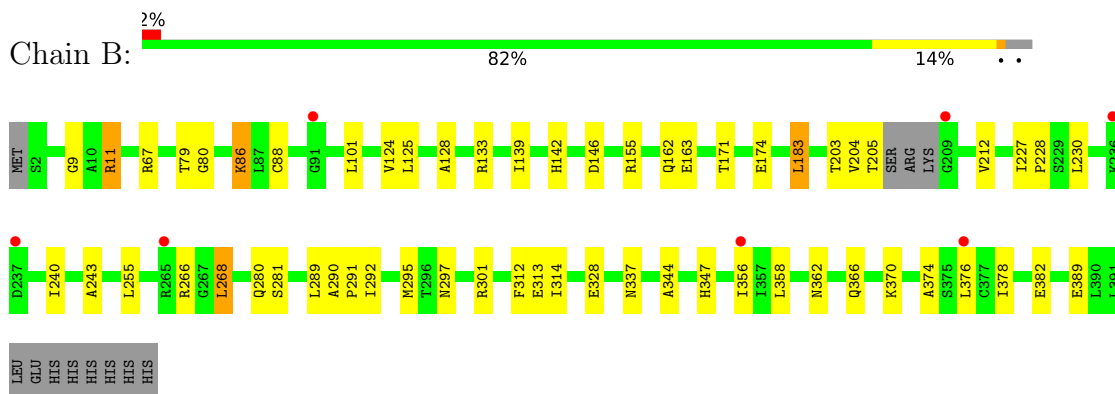
### 3 Residue-property plots [i](#)

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.

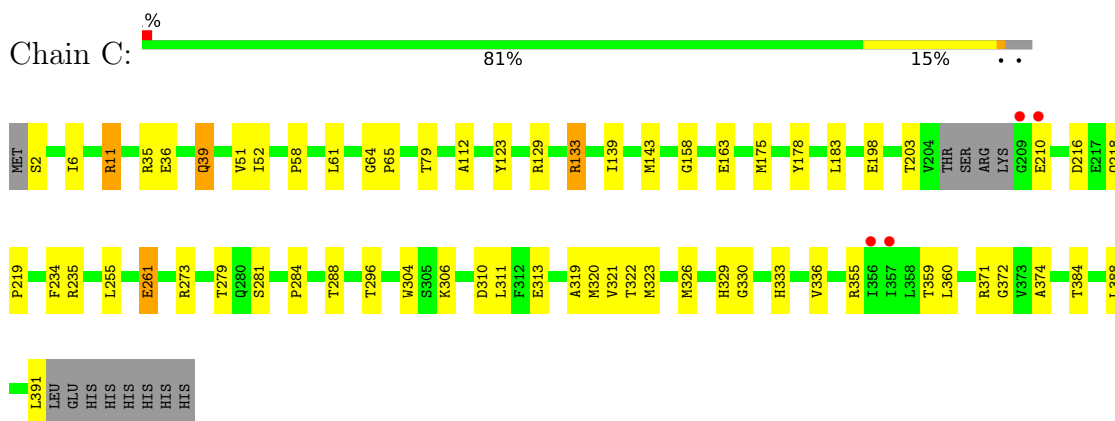
- Molecule 1: Thiolase



- Molecule 1: Thiolase



- Molecule 1: Thiolase



- Molecule 1: Thiolase



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.10Å 93.75Å 111.53Å 90.00° 108.29° 90.00°	Depositor
Resolution (Å)	34.47 – 2.30 34.45 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (34.47-2.30) 97.9 (34.45-2.30)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.83 (at 2.29Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.191 , 0.265 0.198 , 0.266	Depositor DCC
$R_{free}$ test set	3092 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.3	Xtrriage
Anisotropy	0.076	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 35.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11523	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.81	0/2855	0.94	0/3854
1	B	0.78	0/2828	0.94	2/3818 (0.1%)
1	C	0.80	0/2821	0.98	2/3808 (0.1%)
1	D	0.76	0/2821	0.94	0/3808
All	All	0.79	0/11325	0.95	4/15288 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	11	ARG	NE-CZ-NH2	9.23	124.92	120.30
1	C	11	ARG	NE-CZ-NH1	-9.19	115.70	120.30
1	B	11	ARG	CG-CD-NE	5.62	123.59	111.80
1	B	155	ARG	NE-CZ-NH2	-5.25	117.67	120.30

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	2822	0	2869	24	0
1	B	2796	0	2837	35	0
1	C	2789	0	2830	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2789	0	2830	35	0
2	C	6	0	8	2	0
3	A	100	0	0	3	0
3	B	84	0	0	0	0
3	C	78	0	0	2	0
3	D	59	0	0	1	0
All	All	11523	0	11374	131	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 6.

All (131) 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:C:326:MET:HE1	1:C:333:HIS:HA	1.59	0.83
1:C:58:PRO:HA	1:C:61:LEU:HD12	1.63	0.77
1:C:326:MET:CE	1:C:333:HIS:HA	2.15	0.76
2:C:401:GOL:O1	1:D:63:GLN:HB3	1.89	0.73
1:C:175:MET:HA	1:C:323:MET:HE1	1.71	0.72
1:C:39:GLN:OE1	1:C:39:GLN:C	2.29	0.70
1:C:296:THR:HG22	1:C:329:HIS:CE1	2.29	0.68
1:B:313:GLU:HG2	1:B:356:ILE:HB	1.77	0.65
1:A:364:LEU:HB3	1:A:390:LEU:HD11	1.79	0.65
1:D:107:ASN:HD22	1:D:110:MET:HE1	1.60	0.65
1:A:180:ILE:HG23	1:A:224:LEU:HD21	1.77	0.65
1:C:273:ARG:HG2	1:C:391:LEU:HD21	1.79	0.65
1:C:360:LEU:HD22	1:C:388:LEU:HD13	1.81	0.63
1:B:370:LYS:O	1:B:389:GLU:HA	1.99	0.62
1:A:269:LYS:O	1:A:269:LYS:HG3	2.00	0.62
1:A:210:GLU:HG2	3:A:459:HOH:O	2.01	0.61
1:D:101:LEU:HB3	1:D:110:MET:HE1	1.83	0.60
1:C:279:THR:HG22	1:C:384:THR:HG22	1.82	0.60
1:B:314:ILE:H	1:B:337:ASN:HD21	1.50	0.60
1:B:142:HIS:HD2	1:B:146:ASP:OD2	1.86	0.59
1:D:362:ASN:HD21	1:D:366:GLN:NE2	1.99	0.59
1:D:142:HIS:HD2	1:D:146:ASP:OD2	1.85	0.59
1:B:86:LYS:HD3	1:B:378:ILE:HD13	1.86	0.58
1:C:326:MET:HE3	1:C:336:VAL:CB	2.33	0.58
1:D:233:ALA:HB3	1:D:234:PHE:CD2	2.38	0.58
1:B:362:ASN:HD21	1:B:366:GLN:NE2	2.01	0.58
1:C:360:LEU:C	1:C:360:LEU:HD23	2.24	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:326:MET:HE3	1:C:336:VAL:HB	1.85	0.57
1:A:162:GLN:HE22	1:A:240:ILE:H	1.52	0.57
1:B:227:ILE:HB	1:B:228:PRO:HD3	1.86	0.57
1:C:288:THR:HB	1:C:321:VAL:HG22	1.87	0.56
1:B:9:GLY:HA3	1:B:358:LEU:HD21	1.87	0.56
1:C:296:THR:HG21	3:C:506:HOH:O	2.05	0.56
1:B:290:ALA:N	1:B:291:PRO:CD	2.69	0.55
1:D:107:ASN:HD22	1:D:110:MET:CE	2.19	0.55
1:B:67:ARG:NH1	1:B:79:THR:O	2.39	0.55
1:D:36:GLU:HG2	1:D:200:VAL:HB	1.88	0.54
1:C:11:ARG:NH1	1:C:216:ASP:OD1	2.39	0.54
1:C:319:ALA:C	1:C:323:MET:HE2	2.28	0.54
1:D:100:ASP:OD1	1:D:277:HIS:HE1	1.91	0.53
1:B:142:HIS:CD2	1:B:146:ASP:OD2	2.62	0.53
1:A:269:LYS:O	1:A:269:LYS:CG	2.56	0.53
1:D:102:LEU:HG	1:D:110:MET:HE3	1.91	0.52
1:B:125:LEU:HD23	1:B:128:ALA:HB2	1.92	0.52
1:C:52:ILE:O	1:C:112:ALA:HA	2.09	0.51
1:C:326:MET:HE3	1:C:336:VAL:CG1	2.40	0.51
1:C:326:MET:CE	1:C:336:VAL:HB	2.39	0.51
3:C:564:HOH:O	1:D:103:LYS:HE3	2.10	0.51
1:B:203:THR:HG22	1:B:212:VAL:HG22	1.93	0.51
1:C:326:MET:HE3	1:C:336:VAL:HG11	1.92	0.51
1:B:183:LEU:HD22	1:B:344:ALA:HB2	1.93	0.50
1:B:313:GLU:O	1:B:374:ALA:HA	2.11	0.50
1:D:88:CYS:O	1:D:376:LEU:HD11	2.11	0.50
1:D:101:LEU:HB3	1:D:110:MET:CE	2.42	0.50
1:C:178:TYR:OH	1:C:322:THR:HG21	2.12	0.50
1:C:36:GLU:HA	1:C:36:GLU:OE1	2.12	0.50
1:C:51:VAL:HG23	1:C:79:THR:CG2	2.42	0.49
1:D:142:HIS:CD2	1:D:146:ASP:HB2	2.46	0.49
1:B:289:LEU:C	1:B:291:PRO:HD2	2.33	0.49
1:D:58:PRO:HA	1:D:61:LEU:HD12	1.94	0.49
1:D:277:HIS:HD2	3:D:453:HOH:O	1.94	0.49
1:A:124:VAL:HG12	1:B:124:VAL:HG12	1.95	0.48
1:C:311:LEU:O	1:C:372:GLY:HA2	2.14	0.48
2:C:401:GOL:O1	1:D:63:GLN:OE1	2.30	0.48
1:B:313:GLU:HA	1:B:337:ASN:ND2	2.28	0.48
1:C:163:GLU:OE2	1:C:235:ARG:NH2	2.47	0.47
1:D:70:ALA:O	1:D:73:ALA:HB3	2.15	0.47
1:C:158:GLY:HA3	1:C:234:PHE:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:GLU:OE1	1:C:359:THR:OG1	2.26	0.47
1:D:310:ASP:O	1:D:335:LYS:HD2	2.15	0.47
1:A:100:ASP:OD2	1:B:101:LEU:HG	2.15	0.47
1:C:296:THR:HG22	1:C:329:HIS:HE1	1.77	0.47
1:D:312:PHE:HB2	1:D:336:VAL:HG22	1.96	0.47
1:B:88:CYS:O	1:B:376:LEU:HD11	2.15	0.46
1:A:78:ALA:HA	1:B:281:SER:O	2.15	0.46
1:A:158:GLY:HA3	1:A:234:PHE:CZ	2.50	0.46
1:A:361:ILE:O	1:A:365:ARG:HG3	2.16	0.46
1:D:237:ASP:OD2	1:D:237:ASP:N	2.48	0.46
1:A:178:TYR:C	1:A:178:TYR:CD1	2.89	0.46
1:B:171:THR:OG1	1:B:174:GLU:HG3	2.15	0.46
1:B:162:GLN:HE22	1:B:240:ILE:H	1.64	0.46
1:B:255:LEU:HD21	1:B:358:LEU:HG	1.98	0.46
1:B:266:ARG:HB2	1:B:268:LEU:HD22	1.97	0.46
1:B:230:LEU:CD1	1:B:243:ALA:HB2	2.46	0.46
1:D:142:HIS:CD2	1:D:146:ASP:OD2	2.67	0.46
1:B:204:VAL:O	1:B:205:THR:HB	2.16	0.45
1:B:295:MET:CE	1:B:312:PHE:CE1	3.00	0.44
1:C:304:TRP:CZ2	1:C:371:ARG:HD3	2.52	0.44
1:A:133:ARG:O	1:C:129:ARG:HG3	2.16	0.44
1:C:11:ARG:HD3	1:C:198:GLU:HG3	1.99	0.44
1:C:64:GLY:O	1:C:65:PRO:C	2.55	0.44
1:C:310:ASP:O	1:C:311:LEU:HD23	2.18	0.44
1:C:6:ILE:HG21	1:C:255:LEU:HD23	1.99	0.44
1:D:125:LEU:HD23	1:D:128:ALA:HB2	1.99	0.44
1:D:195:LEU:HD21	1:D:339:TYR:HB3	1.99	0.44
1:C:320:MET:HA	1:C:323:MET:HE3	1.99	0.44
1:C:281:SER:O	1:D:78:ALA:HA	2.18	0.44
1:A:80:GLY:HA2	1:B:382:GLU:OE2	2.18	0.44
1:C:218:GLN:N	1:C:219:PRO:CD	2.81	0.43
1:D:102:LEU:HG	1:D:110:MET:CE	2.48	0.43
1:C:123:TYR:HB3	1:C:139:ILE:HG12	2.00	0.43
1:A:125:LEU:N	1:A:125:LEU:HD12	2.34	0.43
1:A:144:PHE:O	1:A:149:GLU:HG2	2.18	0.43
1:D:174:GLU:HB3	1:D:323:MET:HE1	2.01	0.43
1:C:261:GLU:OE2	1:C:261:GLU:N	2.49	0.43
1:D:249:SER:OG	1:D:347:HIS:N	2.52	0.43
1:D:388:LEU:C	1:D:388:LEU:HD12	2.40	0.43
1:A:382:GLU:OE2	1:B:80:GLY:HA2	2.18	0.42
1:A:262:GLU:CD	1:A:265:ARG:HH21	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:THR:CG2	1:C:329:HIS:CE1	3.01	0.42
1:A:124:VAL:HG22	1:A:140:LYS:HB2	2.01	0.42
1:B:86:LYS:HD3	1:B:378:ILE:CD1	2.49	0.42
1:C:313:GLU:O	1:C:374:ALA:HA	2.20	0.42
1:A:2:SER:O	1:A:103:LYS:NZ	2.51	0.41
1:A:97:LEU:HD23	1:B:101:LEU:HD11	2.02	0.41
1:C:6:ILE:CG2	1:C:255:LEU:HD23	2.50	0.41
1:D:125:LEU:CD2	1:D:128:ALA:HB2	2.50	0.41
1:B:292:ILE:HD13	1:B:328:GLU:HG3	2.01	0.41
1:D:218:GLN:N	1:D:219:PRO:CD	2.84	0.41
1:D:173:GLU:CD	1:D:173:GLU:H	2.23	0.41
1:C:39:GLN:OE1	1:C:39:GLN:O	2.38	0.41
1:C:355:ARG:HD2	1:C:355:ARG:C	2.41	0.41
1:D:352:THR:O	1:D:356:ILE:HG12	2.21	0.41
1:B:11:ARG:HH11	1:B:11:ARG:HD2	1.76	0.40
1:D:107:ASN:HB2	1:D:110:MET:HE2	2.03	0.40
1:B:280:GLN:NE2	1:B:297:ASN:HD21	2.19	0.40
1:C:306:LYS:HE3	1:C:330:GLY:O	2.22	0.40
1:D:46:GLU:HA	1:D:76:PRO:HG3	2.03	0.40
1:A:13:PRO:HD3	1:A:200:VAL:O	2.21	0.40
1:A:162:GLN:NE2	3:A:417:HOH:O	2.54	0.40
1:A:277:HIS:HD2	3:A:471:HOH:O	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	388/399 (97%)	374 (96%)	14 (4%)	0	100	100
1	B	383/399 (96%)	371 (97%)	11 (3%)	1 (0%)	41	50
1	C	382/399 (96%)	363 (95%)	18 (5%)	1 (0%)	41	50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	382/399 (96%)	365 (96%)	17 (4%)	0	100	100
All	All	1535/1596 (96%)	1473 (96%)	60 (4%)	2 (0%)	51	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	133	ARG
1	C	133	ARG

### 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	286/295 (97%)	280 (98%)	6 (2%)	53	70
1	B	283/295 (96%)	276 (98%)	7 (2%)	47	65
1	C	282/295 (96%)	272 (96%)	10 (4%)	36	50
1	D	282/295 (96%)	271 (96%)	11 (4%)	32	46
All	All	1133/1180 (96%)	1099 (97%)	34 (3%)	41	57

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	206	SER
1	A	214	LYS
1	A	224	LEU
1	A	236	LYS
1	A	301	ARG
1	B	86	LYS
1	B	139	ILE
1	B	163	GLU
1	B	183	LEU
1	B	268	LEU
1	B	301	ARG

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Mol	Chain	Res	Type
1	B	347	HIS
1	C	2	SER
1	C	35	ARG
1	C	39	GLN
1	C	133	ARG
1	C	143	MET
1	C	183	LEU
1	C	203	THR
1	C	210	GLU
1	C	261	GLU
1	C	284	PRO
1	D	2	SER
1	D	58	PRO
1	D	88	CYS
1	D	133	ARG
1	D	139	ILE
1	D	182	SER
1	D	220	LEU
1	D	229	SER
1	D	264	GLN
1	D	306	LYS
1	D	389	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	162	GLN
1	A	277	HIS
1	A	280	GLN
1	A	347	HIS
1	A	362	ASN
1	A	366	GLN
1	B	68	GLN
1	B	142	HIS
1	B	162	GLN
1	B	218	GLN
1	B	277	HIS
1	B	280	GLN
1	B	337	ASN
1	B	362	ASN
1	C	68	GLN

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Mol	Chain	Res	Type
1	C	218	GLN
1	C	223	ASN
1	C	264	GLN
1	C	280	GLN
1	C	282	GLN
1	C	315	ASN
1	C	362	ASN
1	D	68	GLN
1	D	99	HIS
1	D	107	ASN
1	D	108	GLN
1	D	142	HIS
1	D	277	HIS
1	D	280	GLN
1	D	282	GLN
1	D	315	ASN
1	D	362	ASN
1	D	366	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](#)

1 ligand is 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	C	401	-	5,5,5	0.19	0	5,5,5	0.45	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	C	401	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	401	GOL	O1-C1-C2-C3
2	C	401	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	GOL	2	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	390/399 (97%)	0.07	7 (1%) 68 74	16, 26, 40, 69	0
1	B	387/399 (96%)	0.14	7 (1%) 68 74	18, 28, 47, 73	0
1	C	386/399 (96%)	0.18	4 (1%) 82 86	19, 30, 45, 72	0
1	D	386/399 (96%)	0.18	8 (2%) 63 70	19, 30, 47, 69	0
All	All	1549/1596 (97%)	0.14	26 (1%) 70 76	16, 28, 46, 73	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	209	GLY	4.5
1	C	209	GLY	4.1
1	B	209	GLY	3.4
1	A	237	ASP	3.4
1	A	267	GLY	3.3
1	A	236	LYS	3.1
1	A	207	ARG	2.8
1	B	265	ARG	2.6
1	D	225	GLU	2.5
1	B	376	LEU	2.5
1	B	356	ILE	2.4
1	D	211	SER	2.4
1	A	87	LEU	2.4
1	C	356	ILE	2.3
1	B	91	GLY	2.3
1	B	237	ASP	2.3
1	C	210	GLU	2.3
1	D	348	PRO	2.1
1	B	236	LYS	2.1
1	D	349	VAL	2.1
1	A	349	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	351	SER	2.1
1	A	353	GLY	2.0
1	D	169	GLY	2.0
1	C	357	ILE	2.0
1	D	56	VAL	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(Å <sup>2</sup> )	Q<0.9
2	GOL	C	401	6/6	0.92	0.18	30,34,35,36	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.