



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:36 AM GMT

PDB ID : 2HUF  
Title : Crystal structure of Aedes aegypti alanine glyoxylate aminotransferase  
Authors : Han, Q.; Robinson, H.; Gao, Y.G.; Vogelaar, N.; Wilson, S.R.; Rizzi, M.; Li, J.  
Deposited on : 2006-07-26  
Resolution : 1.75 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.

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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 : 1.7 (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

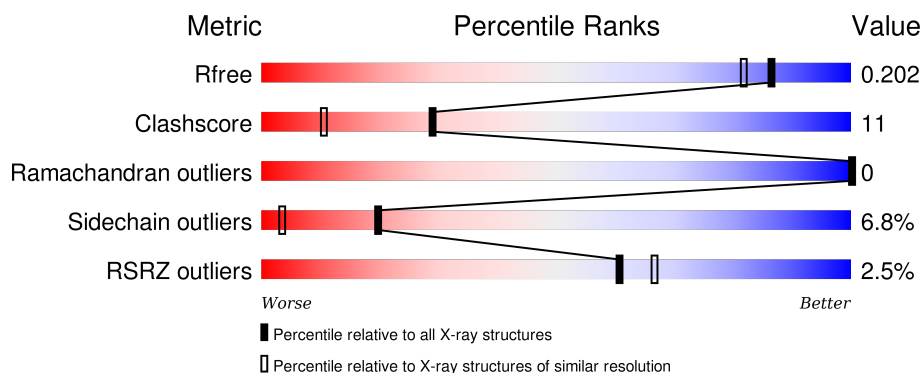
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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.

Mol	Chain	Length	Quality of chain
1	A	393	 3% 79% 17% ...
1	B	393	 2% 79% 16% ...

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	1BO	A	602	-	-	-	X
2	1BO	B	601	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6486 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.

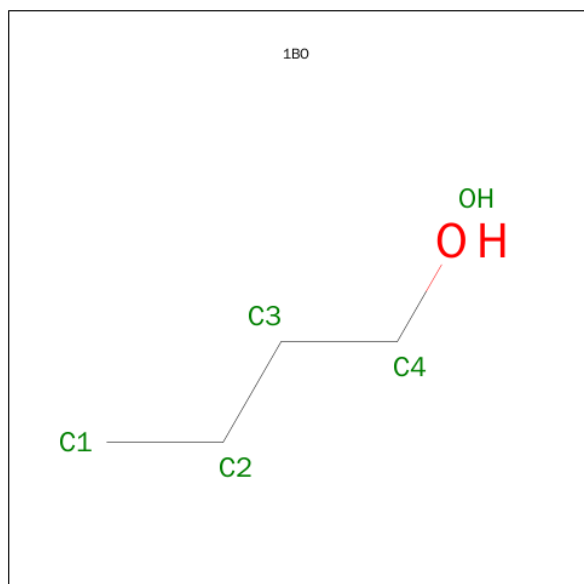
- Molecule 1 is a protein called Alanine glyoxylate aminotransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	385	Total	C	N	O	P	S	0	0	0
			2978	1885	523	549	1	20			
1	B	385	Total	C	N	O	P	S	0	0	0
			2978	1885	523	549	1	20			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	206	LLP	LYS	MODIFIED RESIDUE	UNP Q3LSM4
B	206	LLP	LYS	MODIFIED RESIDUE	UNP Q3LSM4

- Molecule 2 is 1-BUTANOL (three-letter code: 1BO) (formula: C<sub>4</sub>H<sub>10</sub>O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			5	4	1		

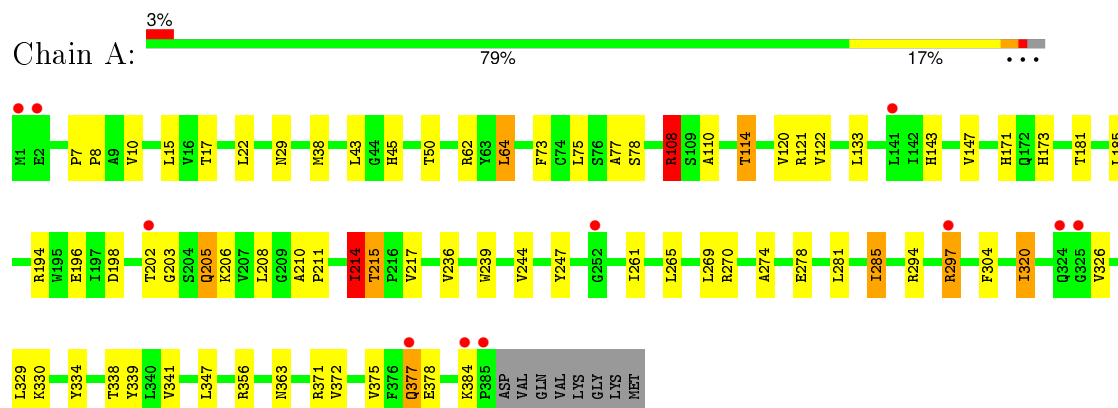
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	255	Total	O	0	0
			255	255		
3	B	265	Total	O	0	0
			265	265		

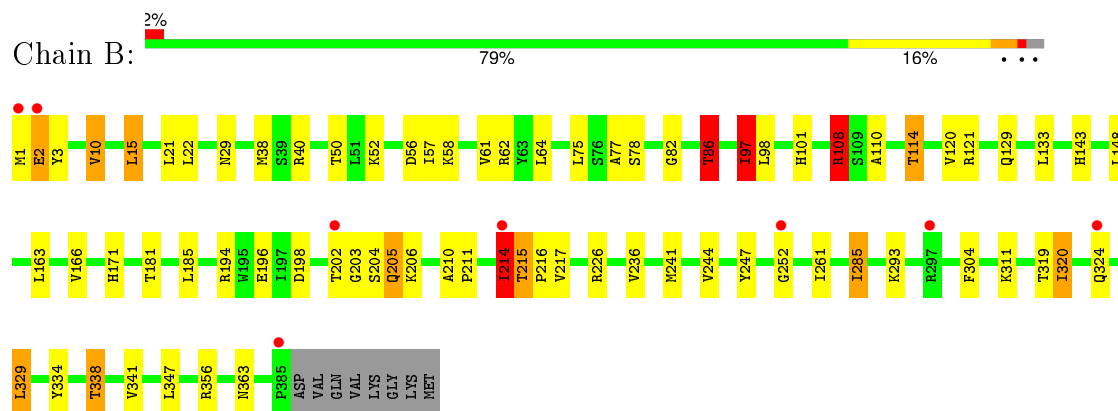
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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: Alanine glyoxylate aminotransferase



- Molecule 1: Alanine glyoxylate aminotransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	136.96Å 136.96Å 120.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.20 – 1.75 29.19 – 1.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.20-1.75) 99.9 (29.19-1.75)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.19 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.172 , 0.203 0.171 , 0.202	Depositor DCC
$R_{free}$ test set	4233 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.8	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 44.0	EDS
Estimated twinning fraction	0.017 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 84776 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6486	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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, 0.375 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: LLP, 1BO

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	0.60	0/3017	0.85	9/4092 (0.2%)
1	B	0.61	0/3017	0.87	9/4092 (0.2%)
All	All	0.61	0/6034	0.86	18/8184 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	ARG	NE-CZ-NH2	-14.35	113.12	120.30
1	A	108	ARG	NE-CZ-NH2	-14.23	113.19	120.30
1	B	108	ARG	NE-CZ-NH1	13.67	127.13	120.30
1	A	108	ARG	NE-CZ-NH1	12.39	126.50	120.30
1	A	62	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	B	97	ILE	CB-CA-C	-7.55	96.50	111.60
1	A	62	ARG	NE-CZ-NH1	7.54	124.07	120.30
1	B	214	ILE	CB-CA-C	-6.93	97.75	111.60
1	B	62	ARG	NE-CZ-NH2	-6.70	116.95	120.30
1	B	62	ARG	NE-CZ-NH1	6.65	123.63	120.30
1	B	10	VAL	CG1-CB-CG2	6.16	120.75	110.90
1	A	108	ARG	CD-NE-CZ	6.06	132.08	123.60
1	A	214	ILE	CB-CA-C	-6.02	99.55	111.60
1	B	108	ARG	CD-NE-CZ	5.92	131.89	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	239	TRP	CA-CB-CG	5.41	123.97	113.70
1	A	64	LEU	CB-CG-CD1	5.34	120.07	111.00
1	A	214	ILE	CG1-CB-CG2	5.21	122.86	111.40
1	B	86	THR	N-CA-CB	-5.10	100.61	110.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	252	GLY	Peptide

## 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	2978	0	2987	77	0
1	B	2978	0	2987	67	0
2	A	5	0	10	2	0
2	B	5	0	10	3	0
3	A	255	0	0	12	0
3	B	265	0	0	11	0
All	All	6486	0	5994	136	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 11.

All (136) 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:108:ARG:HD2	3:A:712:HOH:O	1.62	0.99
1:B:108:ARG:HD2	3:B:862:HOH:O	1.65	0.94
1:B:22:LEU:H	1:B:29:ASN:HD21	1.15	0.91
1:A:22:LEU:H	1:A:29:ASN:HD21	1.17	0.88
1:B:293:LYS:HD2	3:B:765:HOH:O	1.73	0.87
1:B:334:TYR:O	1:B:338:THR:HG23	1.79	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:ARG:O	1:A:297:ARG:HD3	1.80	0.81
1:B:75:LEU:CD1	1:B:217:VAL:HG21	2.11	0.81
1:A:45:HIS:NE2	2:B:601:1BO:H12	1.97	0.80
1:B:75:LEU:HD12	1:B:217:VAL:HG21	1.64	0.79
1:A:121:ARG:HH11	1:A:143:HIS:HD2	1.28	0.78
1:A:371:ARG:O	1:A:375:VAL:HG12	1.83	0.77
1:B:205:GLN:H	1:B:205:GLN:HE21	1.33	0.76
1:B:202:THR:HG22	1:B:203:GLY:H	1.50	0.76
1:A:114:THR:HG22	1:A:120:VAL:HG21	1.68	0.76
1:B:114:THR:HG22	1:B:120:VAL:HG21	1.68	0.75
1:A:247:TYR:OH	1:A:261:ILE:HG13	1.87	0.74
1:A:304:PHE:HB3	1:A:320:ILE:HD13	1.69	0.74
1:A:181:THR:O	1:A:202:THR:HG23	1.88	0.74
1:A:202:THR:HG22	1:A:203:GLY:H	1.50	0.74
1:A:10:VAL:HG21	1:B:56:ASP:OD1	1.89	0.71
1:B:171:HIS:HE1	1:B:198:ASP:OD2	1.74	0.71
1:A:304:PHE:HB3	1:A:320:ILE:CD1	2.20	0.70
1:A:110:ALA:O	1:A:114:THR:HG23	1.93	0.69
1:A:205:GLN:HE21	1:A:205:GLN:H	1.41	0.69
1:B:247:TYR:OH	1:B:261:ILE:HG13	1.93	0.68
1:A:334:TYR:O	1:A:338:THR:HG23	1.94	0.67
1:B:77:ALA:O	1:B:215:THR:HG23	1.95	0.66
1:A:75:LEU:CD1	1:A:217:VAL:HG11	2.25	0.66
1:A:171:HIS:HE1	1:A:198:ASP:OD2	1.77	0.66
1:A:77:ALA:O	1:A:215:THR:HG23	1.95	0.66
1:B:181:THR:O	1:B:202:THR:HG23	1.96	0.65
1:B:163:LEU:O	1:B:166:VAL:HG22	1.96	0.65
1:A:210:ALA:HB1	1:A:211:PRO:HD2	1.77	0.64
1:A:29:ASN:HD22	1:A:363:ASN:HD21	1.45	0.64
1:A:194:ARG:HH11	1:A:194:ARG:HG3	1.63	0.64
1:B:108:ARG:CD	3:B:862:HOH:O	2.33	0.63
1:A:10:VAL:HG21	1:B:52:LYS:HZ2	1.62	0.63
1:B:205:GLN:N	1:B:205:GLN:HE21	1.97	0.62
1:B:121:ARG:HH11	1:B:143:HIS:HD2	1.44	0.62
1:A:121:ARG:HH11	1:A:143:HIS:CD2	2.14	0.62
1:B:110:ALA:O	1:B:114:THR:HG23	1.99	0.62
1:B:97:ILE:CD1	1:B:120:VAL:HG22	2.31	0.60
1:B:356:ARG:HH12	2:B:601:1BO:H42	1.68	0.59
1:A:202:THR:HG22	1:A:203:GLY:N	2.17	0.59
1:A:17:THR:OG1	1:B:40:ARG:NE	2.35	0.59
1:B:50:THR:HG21	1:B:261:ILE:HD11	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:LEU:HD12	1:A:217:VAL:HG11	1.86	0.57
1:B:97:ILE:HD12	1:B:120:VAL:HA	1.85	0.57
1:B:29:ASN:HD22	1:B:363:ASN:HD21	1.52	0.57
1:A:205:GLN:HE21	1:A:205:GLN:N	2.02	0.57
1:B:215:THR:HG22	3:B:608:HOH:O	2.05	0.56
1:A:281:LEU:O	1:A:285:ILE:HG23	2.07	0.55
1:A:206:LLP:O3	1:A:206:LLP:NZ	2.37	0.55
1:B:206:LLP:NZ	1:B:206:LLP:O3	2.38	0.55
1:A:50:THR:HG21	1:A:261:ILE:HD11	1.88	0.55
1:B:82:GLY:O	1:B:86:THR:HB	2.07	0.55
1:B:202:THR:HG22	1:B:203:GLY:N	2.18	0.54
1:A:10:VAL:HG23	3:B:747:HOH:O	2.07	0.54
1:B:97:ILE:HD11	1:B:120:VAL:HG22	1.88	0.54
1:B:210:ALA:HB1	1:B:211:PRO:HD2	1.91	0.53
1:B:205:GLN:H	1:B:205:GLN:NE2	2.03	0.53
1:A:247:TYR:OH	1:A:261:ILE:CG1	2.55	0.53
1:A:10:VAL:CG2	1:B:52:LYS:NZ	2.72	0.53
1:B:143:HIS:HE1	3:B:716:HOH:O	1.90	0.53
1:A:143:HIS:HE1	3:A:817:HOH:O	1.90	0.52
1:A:10:VAL:HG21	1:B:52:LYS:NZ	2.24	0.52
1:A:171:HIS:HD2	1:A:196:GLU:OE2	1.92	0.52
1:A:270:ARG:HD2	1:B:15:LEU:HG	1.92	0.51
1:A:108:ARG:CD	3:A:712:HOH:O	2.36	0.51
1:A:294:ARG:O	1:A:297:ARG:CD	2.55	0.51
1:A:77:ALA:O	1:A:215:THR:CG2	2.58	0.51
1:A:341:VAL:HG21	1:A:375:VAL:HG11	1.93	0.50
1:A:38:MET:SD	1:A:211:PRO:HG3	2.52	0.50
1:B:75:LEU:HD11	1:B:217:VAL:HG21	1.94	0.50
1:B:247:TYR:OH	1:B:261:ILE:CG1	2.59	0.50
1:A:377:GLN:HG2	1:A:378:GLU:N	2.22	0.50
1:A:274:ALA:O	1:A:278:GLU:HG3	2.12	0.49
1:A:356:ARG:HH22	2:A:602:1BO:H31	1.77	0.49
1:A:377:GLN:CG	3:A:775:HOH:O	2.61	0.48
1:A:384:LYS:HE2	3:A:790:HOH:O	2.12	0.48
1:A:29:ASN:ND2	1:A:363:ASN:HD21	2.11	0.48
1:A:22:LEU:H	1:A:29:ASN:ND2	1.99	0.48
1:A:377:GLN:HG2	3:A:775:HOH:O	2.13	0.48
1:A:214:ILE:HG12	1:A:265:LEU:HB3	1.95	0.47
1:A:75:LEU:HD11	1:A:217:VAL:HG11	1.93	0.47
1:B:194:ARG:NE	3:B:841:HOH:O	2.47	0.47
1:B:285:ILE:HD13	3:B:733:HOH:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:SER:CB	1:B:214:ILE:HD12	2.45	0.47
1:B:57:ILE:O	1:B:61:VAL:HG12	2.15	0.47
1:B:75:LEU:HD12	1:B:217:VAL:CG2	2.38	0.46
1:B:129:GLN:HE21	1:B:311:LYS:NZ	2.12	0.46
1:A:304:PHE:HB3	1:A:320:ILE:HD11	1.97	0.46
1:A:78:SER:C	1:A:215:THR:HG21	2.36	0.46
1:B:171:HIS:HD2	1:B:196:GLU:OE2	1.99	0.46
1:B:261:ILE:HD12	1:B:261:ILE:O	2.17	0.45
1:B:101:HIS:HD2	3:B:696:HOH:O	1.98	0.45
1:A:205:GLN:H	1:A:205:GLN:NE2	2.10	0.45
1:B:40:ARG:NE	3:B:715:HOH:O	2.47	0.45
1:A:114:THR:HG21	3:A:758:HOH:O	2.17	0.45
1:A:356:ARG:HH22	2:A:602:1BO:C3	2.30	0.45
1:B:319:THR:HG22	1:B:356:ARG:HG2	1.99	0.44
1:B:215:THR:HA	1:B:216:PRO:HD3	1.83	0.44
1:A:78:SER:O	1:A:215:THR:HG21	2.17	0.44
1:A:110:ALA:O	1:A:114:THR:CG2	2.64	0.44
1:A:73:PHE:CE1	1:A:244:VAL:HG13	2.52	0.44
1:A:261:ILE:O	1:A:261:ILE:HD12	2.17	0.44
1:B:356:ARG:HH12	2:B:601:1BO:C4	2.29	0.44
1:A:185:LEU:HB3	1:A:202:THR:HG21	1.98	0.44
1:B:78:SER:C	1:B:215:THR:HG21	2.37	0.43
1:A:372:VAL:HA	1:A:375:VAL:CG1	2.48	0.43
1:A:10:VAL:HG23	1:B:52:LYS:HZ1	1.83	0.43
1:B:2:GLU:HG3	1:B:3:TYR:N	2.33	0.43
1:B:185:LEU:HB3	1:B:202:THR:HG21	1.99	0.43
1:A:261:ILE:HD12	3:A:749:HOH:O	2.18	0.43
1:B:58:LYS:HA	1:B:61:VAL:HG12	2.01	0.43
1:B:58:LYS:HA	1:B:61:VAL:CG1	2.49	0.43
1:B:171:HIS:CE1	1:B:198:ASP:OD2	2.64	0.42
1:A:215:THR:HG22	3:A:621:HOH:O	2.18	0.42
1:A:173:HIS:HE1	3:A:780:HOH:O	2.01	0.42
1:A:73:PHE:CE1	1:A:244:VAL:CG1	3.02	0.42
1:A:7:PRO:HA	1:A:8:PRO:HD3	1.91	0.42
1:A:214:ILE:CG1	1:A:265:LEU:HB3	2.50	0.42
1:A:214:ILE:HD13	1:A:269:LEU:HB2	2.01	0.42
1:A:122:VAL:CG1	3:A:785:HOH:O	2.67	0.42
1:B:110:ALA:O	1:B:114:THR:CG2	2.67	0.42
1:B:329:LEU:HD22	3:B:845:HOH:O	2.19	0.42
1:A:10:VAL:CG2	1:B:56:ASP:OD1	2.65	0.42
1:A:45:HIS:H	1:A:45:HIS:HD1	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:TYR:HB3	1:A:341:VAL:HG13	2.02	0.41
1:B:98:LEU:HD23	1:B:148:LEU:HD13	2.02	0.41
1:B:29:ASN:ND2	1:B:363:ASN:HD21	2.16	0.41
1:B:21:LEU:HB3	1:B:341:VAL:HB	2.02	0.41
1:B:304:PHE:HB3	1:B:320:ILE:HD13	2.02	0.41
1:B:38:MET:SD	1:B:211:PRO:HG3	2.61	0.41
1:A:377:GLN:HB2	3:A:650:HOH:O	2.20	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	382/393 (97%)	378 (99%)	4 (1%)	0	100	100
1	B	382/393 (97%)	377 (99%)	5 (1%)	0	100	100
All	All	764/786 (97%)	755 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	315/322 (98%)	295 (94%)	20 (6%)	22	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	315/322 (98%)	292 (93%)	23 (7%)	17	3
All	All	630/644 (98%)	587 (93%)	43 (7%)	20	3

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	43	LEU
1	A	64	LEU
1	A	108	ARG
1	A	114	THR
1	A	133	LEU
1	A	147	VAL
1	A	205	GLN
1	A	208	LEU
1	A	214	ILE
1	A	215	THR
1	A	236	VAL
1	A	285	ILE
1	A	297	ARG
1	A	320	ILE
1	A	326	VAL
1	A	329	LEU
1	A	330	LYS
1	A	347	LEU
1	A	377	GLN
1	B	1	MET
1	B	2	GLU
1	B	10	VAL
1	B	15	LEU
1	B	64	LEU
1	B	86	THR
1	B	97	ILE
1	B	108	ARG
1	B	114	THR
1	B	133	LEU
1	B	205	GLN
1	B	214	ILE
1	B	215	THR
1	B	226	ARG
1	B	236	VAL
1	B	241	MET

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Mol	Chain	Res	Type
1	B	244	VAL
1	B	285	ILE
1	B	320	ILE
1	B	324	GLN
1	B	329	LEU
1	B	338	THR
1	B	347	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	69	ASN
1	A	101	HIS
1	A	129	GLN
1	A	143	HIS
1	A	171	HIS
1	A	173	HIS
1	A	205	GLN
1	A	333	GLN
1	B	29	ASN
1	B	69	ASN
1	B	101	HIS
1	B	129	GLN
1	B	143	HIS
1	B	152	GLN
1	B	171	HIS
1	B	172	GLN
1	B	205	GLN
1	B	333	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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
1	LLP	A	206	1	23,24,25	1.73	5 (21%)	28,32,34	2.13	8 (28%)
1	LLP	B	206	1	23,24,25	1.76	5 (21%)	28,32,34	2.14	7 (25%)

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
1	LLP	A	206	1	-	0/15/17/19	0/1/1/1
1	LLP	B	206	1	-	0/15/17/19	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	206	LLP	O3-C3	-5.05	1.25	1.37
1	A	206	LLP	O3-C3	-4.90	1.25	1.37
1	A	206	LLP	P-OP3	-2.08	1.47	1.54
1	B	206	LLP	CD-CE	2.08	1.58	1.51
1	A	206	LLP	CD-CE	2.09	1.58	1.51
1	B	206	LLP	C6-N1	2.28	1.39	1.34
1	A	206	LLP	C4'-NZ	2.86	1.36	1.27
1	B	206	LLP	C4'-NZ	2.91	1.36	1.27
1	B	206	LLP	C4-C4'	3.03	1.52	1.46
1	A	206	LLP	C4-C4'	3.10	1.52	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	LLP	C4-C4'-NZ	-3.33	106.54	125.06
1	B	206	LLP	OP2-P-OP4	-3.27	97.15	106.56
1	A	206	LLP	OP2-P-OP4	-3.04	97.80	106.56
1	B	206	LLP	O-C-CA	-3.03	117.61	125.49
1	B	206	LLP	C5'-C5-C6	-2.89	113.82	119.28
1	A	206	LLP	O-C-CA	-2.78	118.25	125.49
1	B	206	LLP	C4-C4'-NZ	-2.72	109.92	125.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	206	LLP	C5'-C5-C6	-2.62	114.33	119.28
1	A	206	LLP	C5-C6-N1	-2.25	119.95	123.86
1	B	206	LLP	C5-C6-N1	-2.11	120.20	123.86
1	A	206	LLP	OP3-P-OP2	2.18	115.67	107.38
1	B	206	LLP	C5'-C5-C4	2.44	125.57	121.47
1	A	206	LLP	C5'-C5-C4	2.50	125.67	121.47
1	A	206	LLP	OP4-C5'-C5	7.28	121.02	108.99
1	B	206	LLP	OP4-C5'-C5	7.65	121.64	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	206	LLP	1	0
1	B	206	LLP	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 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	1BO	A	602	-	4,4,4	0.23	0	3,3,3	0.43	0
2	1BO	B	601	-	4,4,4	0.18	0	3,3,3	0.58	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	1BO	A	602	-	-	0/2/2/2	0/0/0/0
2	1BO	B	601	-	-	0/2/2/2	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.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	602	1BO	2	0
2	B	601	1BO	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 [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	384/393 (97%)	-0.11	11 (2%) 55 61	11, 17, 26, 48	0
1	B	384/393 (97%)	-0.09	8 (2%) 67 73	12, 16, 26, 50	0
All	All	768/786 (97%)	-0.10	19 (2%) 61 67	11, 16, 26, 50	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	8.4
1	A	1	MET	6.6
1	B	385	PRO	5.2
1	B	324	GLN	5.0
1	B	252	GLY	4.5
1	A	141	LEU	4.2
1	A	324	GLN	3.9
1	A	385	PRO	3.6
1	B	2	GLU	3.3
1	A	202	THR	2.9
1	A	377	GLN	2.8
1	A	297	ARG	2.8
1	B	214	ILE	2.5
1	B	202	THR	2.5
1	A	325	GLY	2.4
1	A	384	LYS	2.3
1	B	297	ARG	2.2
1	A	2	GLU	2.2
1	A	252	GLY	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	LLP	B	206	24/25	0.98	0.11	-	13,14,18,20	0
1	LLP	A	206	24/25	0.98	0.12	-	12,15,18,20	0

### 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	1BO	B	601	5/5	0.84	0.21	13.23	28,29,31,32	0
2	1BO	A	602	5/5	0.89	0.16	8.77	30,32,35,36	0

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

There are no such residues in this entry.