



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:45 AM UTC

PDB ID : 6XML / pdb_00006xml
Title : Human aldolase A I98C
Authors : Meneely, K.M.; Brewer, K.; Lamb, A.L.
Deposited on : 2020-06-30
Resolution : 1.88 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

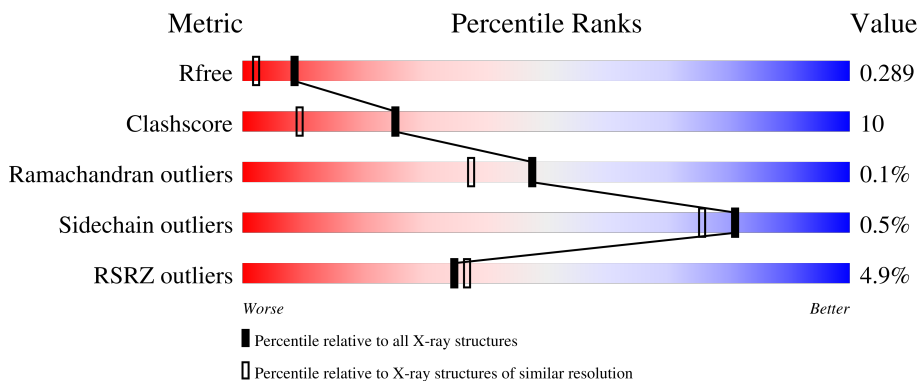
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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}	180053	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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 ≥ 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	364	
1	B	364	

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	PO4	B	402	-	-	X	-
3	GOL	A	405	-	X	-	-
3	GOL	B	405	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11179 atoms, of which 5486 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.

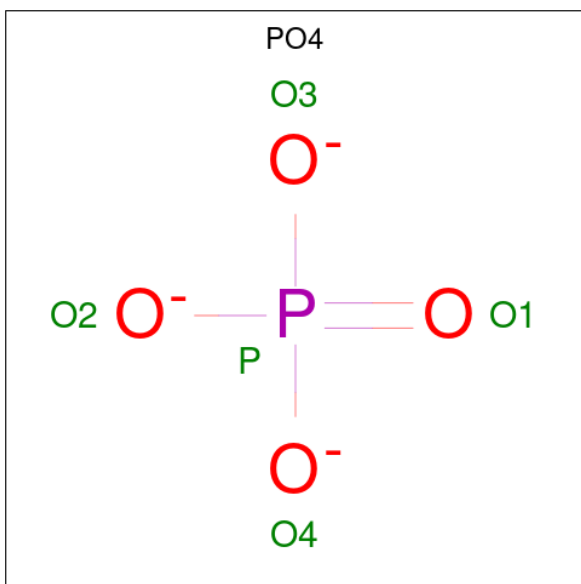
- Molecule 1 is a protein called Fructose-bisphosphate aldolase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	353	5380	1682	2705	472	509	12	0	0	0
1	B	358	5444	1704	2733	478	517	12	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	98	CYS	ILE	engineered mutation	UNP P04075
B	98	CYS	ILE	engineered mutation	UNP P04075

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



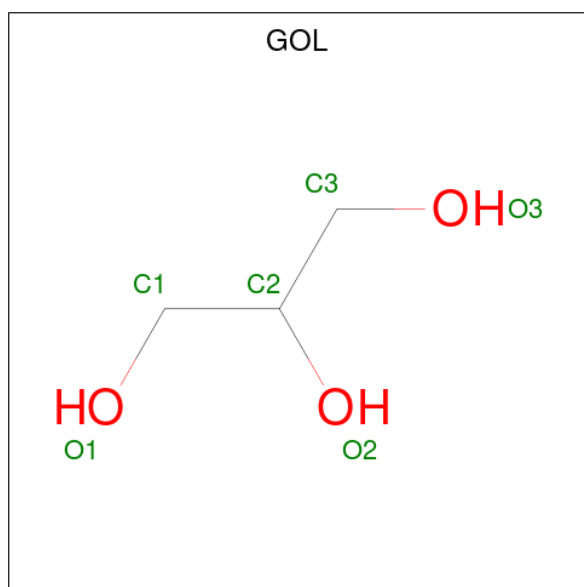
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0

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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		

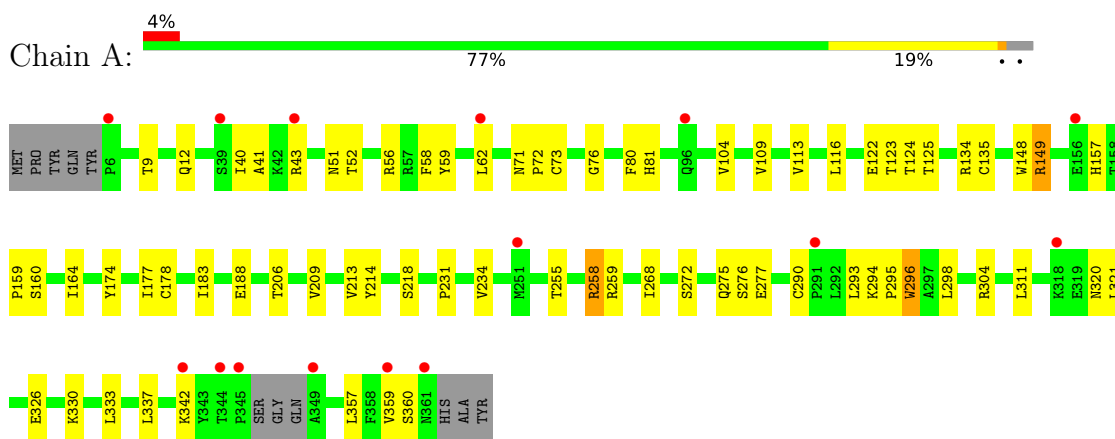
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total 126	O 126	0	0
4	B	125	Total 125	O 125	0	0

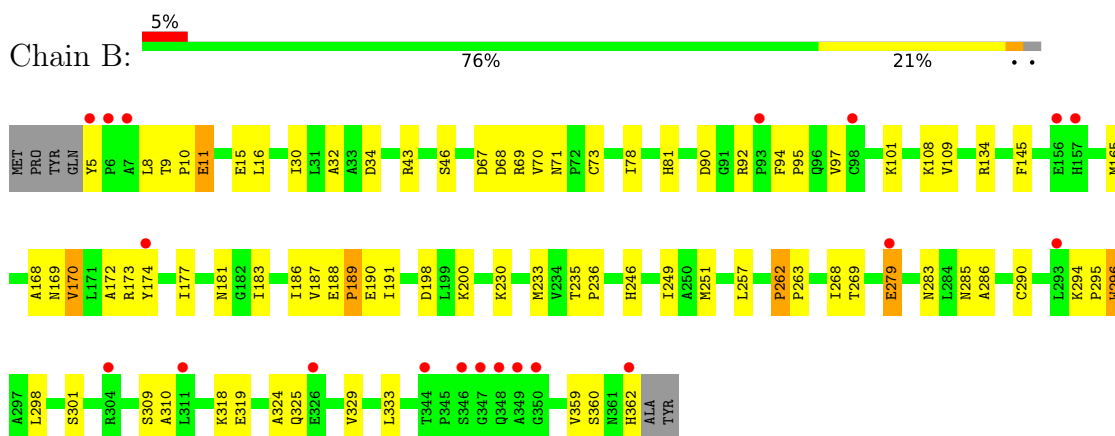
3 Residue-property plots [i](#)

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: Fructose-bisphosphate aldolase A



- Molecule 1: Fructose-bisphosphate aldolase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	162.57Å 162.57Å 167.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.20 – 1.88 40.20 – 1.88	Depositor EDS
% Data completeness (in resolution range)	61.7 (40.20-1.88) 61.7 (40.20-1.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.88Å)	Xtrriage
Refinement program	REFMAC 1.16_3549, PHENIX 1.16_3549	Depositor
R, R_{free}	0.245 , 0.286 (Not available) , 0.289	Depositor DCC
R_{free} test set	3143 reflections (2.97%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtrriage
Anisotropy	0.031	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 32.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11179	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.41 % 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 2.0494e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: PO4, 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	1.02	2/2723 (0.1%)	1.09	4/3687 (0.1%)
1	B	1.15	15/2761 (0.5%)	1.06	5/3741 (0.1%)
All	All	1.08	17/5484 (0.3%)	1.08	9/7428 (0.1%)

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	A	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	187	VAL	C-O	-11.56	1.11	1.24
1	B	10	PRO	C-O	-10.97	1.11	1.24
1	B	188	GLU	C-O	-9.68	1.15	1.24
1	B	9	THR	C-O	-8.86	1.10	1.23
1	B	296	TRP	C-O	-7.83	1.14	1.23
1	A	296	TRP	C-O	-7.49	1.14	1.23
1	B	189	PRO	C-O	-7.39	1.14	1.24
1	B	188	GLU	CD-OE1	-7.12	1.11	1.25
1	B	11	GLU	C-O	-6.78	1.16	1.24
1	B	8	LEU	C-O	-6.55	1.15	1.23
1	B	262	PRO	CA-C	-6.43	1.48	1.51
1	B	191	ILE	C-O	-6.19	1.17	1.24
1	B	190	GLU	C-O	-6.14	1.16	1.24
1	A	231	PRO	N-CD	-6.13	1.39	1.47
1	B	279	GLU	CG-CD	5.98	1.67	1.52
1	B	362	HIS	CA-CB	5.77	1.65	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	186	ILE	C-O	-5.48	1.18	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	TYR	CA-C-N	-7.72	112.75	120.31
1	B	5	TYR	C-N-CA	-7.72	112.75	120.31
1	A	214	TYR	CA-CB-CG	-6.40	102.37	113.90
1	A	258	ARG	CA-C-N	-5.79	111.90	122.38
1	A	258	ARG	C-N-CA	-5.79	111.90	122.38
1	B	170	VAL	N-CA-CB	5.37	120.49	110.77
1	B	5	TYR	O-C-N	-5.36	114.43	123.00
1	A	149	ARG	NE-CZ-NH2	-5.22	114.50	119.20
1	B	262	PRO	O-C-N	-5.07	118.98	121.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	THR	Peptide

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	2675	2705	2704	49	0
1	B	2711	2733	2731	59	0
2	A	10	0	0	1	0
2	B	10	0	0	2	0
3	A	18	24	24	3	0
3	B	18	24	24	7	0
4	A	126	0	0	1	0
4	B	125	0	0	4	0
All	All	5693	5486	5483	106	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 10.

All (106) 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:124:THR:HG22	1:A:125:THR:H	1.00	1.09
1:B:290:CYS:O	1:B:294:LYS:NZ	1.86	1.08
1:A:124:THR:HG22	1:A:125:THR:N	1.80	0.96
1:A:124:THR:CG2	1:A:125:THR:H	1.81	0.93
1:B:319:GLU:OE1	4:B:501:HOH:O	2.01	0.78
1:B:200:LYS:NZ	2:B:402:PO4:O4	2.21	0.74
1:A:304:ARG:HB3	1:A:357:LEU:HD13	1.76	0.68
1:A:149:ARG:HG3	1:A:188:GLU:HG2	1.76	0.67
1:A:272:SER:O	1:A:275:GLN:HG3	1.96	0.66
1:B:230:LYS:HG3	1:B:269:THR:O	1.95	0.66
1:B:68:ASP:OD1	4:B:502:HOH:O	2.14	0.65
1:A:43:ARG:O	1:A:311:LEU:HD21	1.96	0.64
1:A:290:CYS:O	1:A:294:LYS:NZ	2.30	0.64
3:B:405:GOL:HO1	3:B:405:GOL:HO3	1.46	0.64
1:B:108:LYS:HE3	3:B:403:GOL:H32	1.81	0.61
1:A:326:GLU:O	1:A:330:LYS:HG3	2.03	0.59
1:B:359:VAL:O	1:B:360:SER:C	2.46	0.58
1:B:81:HIS:HE2	3:B:405:GOL:H2	1.68	0.58
1:A:337:LEU:HD22	1:A:342:LYS:HD3	1.85	0.58
1:B:165:MET:SD	1:B:165:MET:C	2.86	0.58
1:B:233:MET:HE2	1:B:283:ASN:HB3	1.86	0.58
1:B:67:ASP:O	1:B:70:VAL:HG22	2.05	0.57
1:B:251:MET:HE2	4:B:624:HOH:O	2.05	0.57
1:B:318:LYS:HG3	1:B:318:LYS:O	2.05	0.56
1:A:157:HIS:HA	3:A:404:GOL:H32	1.85	0.56
1:A:304:ARG:HD3	2:A:401:PO4:O1	2.05	0.56
1:A:160:SER:O	1:A:164:ILE:HG13	2.06	0.56
1:B:109:VAL:HG21	1:B:174:TYR:HE1	1.71	0.55
1:A:116:LEU:HD13	1:A:124:THR:OG1	2.07	0.55
1:A:135:CYS:HB3	1:A:183:ILE:HD12	1.90	0.54
1:B:43:ARG:NH1	1:B:43:ARG:HG2	2.23	0.54
1:A:149:ARG:HG3	1:A:188:GLU:CG	2.39	0.53
1:B:90:ASP:OD1	1:B:92:ARG:HG3	2.09	0.53
1:A:41:ALA:HB2	1:A:51:ASN:ND2	2.23	0.53
1:B:257:LEU:HD13	1:B:268:ILE:HD13	1.91	0.52
1:A:80:PHE:O	1:A:81:HIS:C	2.51	0.52
1:A:293:LEU:HD22	1:B:295:PRO:HB3	1.90	0.52
1:B:169:ASN:O	1:B:172:ALA:HB3	2.11	0.51
1:A:209:VAL:O	1:A:213:VAL:HG23	2.11	0.51
1:B:359:VAL:O	1:B:359:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:MET:CE	1:B:283:ASN:HB3	2.41	0.50
1:B:170:VAL:HG12	3:B:404:GOL:H2	1.92	0.50
1:B:43:ARG:HG2	1:B:43:ARG:HH11	1.76	0.49
1:B:168:ALA:O	1:B:169:ASN:C	2.55	0.49
1:B:324:ALA:O	1:B:325:GLN:C	2.56	0.49
1:B:97:VAL:O	1:B:101:LYS:HG3	2.12	0.49
1:B:170:VAL:HG13	3:B:404:GOL:O3	2.12	0.49
1:A:320:ASN:O	1:A:321:LEU:C	2.55	0.49
1:B:109:VAL:O	1:B:134:ARG:NH2	2.42	0.48
1:B:181:ASN:OD1	4:B:503:HOH:O	2.20	0.48
1:B:67:ASP:OD2	1:B:69:ARG:HD3	2.13	0.48
1:B:81:HIS:NE2	3:B:405:GOL:H2	2.28	0.48
1:B:173:ARG:O	1:B:177:ILE:HG13	2.14	0.48
1:B:34:ASP:OD2	1:B:108:LYS:HE2	2.14	0.48
1:B:169:ASN:O	1:B:172:ALA:N	2.45	0.48
1:A:76:GLY:HA2	1:A:104:VAL:O	2.14	0.47
1:A:109:VAL:O	1:A:134:ARG:NH2	2.45	0.47
1:B:30:ILE:HB	1:B:301:SER:HA	1.96	0.47
1:B:43:ARG:O	1:B:46:SER:OG	2.31	0.47
1:A:73:CYS:SG	1:A:333:LEU:HD23	2.55	0.47
1:A:255:THR:O	1:A:259:ARG:HG3	2.15	0.46
1:A:40:ILE:O	1:A:40:ILE:HG13	2.14	0.46
1:A:295:PRO:HG3	1:B:263:PRO:HG3	1.97	0.46
1:B:32:ALA:HB1	1:B:34:ASP:OD1	2.16	0.46
1:A:337:LEU:CD2	1:A:342:LYS:HD3	2.45	0.46
1:B:198:ASP:C	1:B:198:ASP:OD1	2.57	0.46
1:B:294:LYS:HD2	1:B:298:LEU:HD12	1.98	0.45
1:A:157:HIS:HA	3:A:404:GOL:C3	2.46	0.45
1:B:71:ASN:OD1	1:B:101:LYS:HB3	2.16	0.45
1:A:58:PHE:CZ	1:A:321:LEU:HD12	2.52	0.45
1:A:268:ILE:HB	1:A:298:LEU:HD23	1.99	0.45
1:B:235:THR:HB	1:B:236:PRO:HD2	1.98	0.45
1:A:148:TRP:HB2	1:A:174:TYR:CE1	2.53	0.44
1:A:59:TYR:O	1:A:62:LEU:HB3	2.17	0.44
1:B:30:ILE:HG21	1:B:78:ILE:HD11	2.00	0.44
1:A:258:ARG:O	1:B:263:PRO:HD2	2.18	0.44
1:A:218:SER:O	3:A:405:GOL:H12	2.18	0.44
1:A:149:ARG:HG3	1:A:188:GLU:CD	2.42	0.43
1:B:70:VAL:CG1	1:B:329:VAL:HG22	2.48	0.43
1:A:71:ASN:N	1:A:72:PRO:CD	2.82	0.43
1:A:294:LYS:HD2	1:A:298:LEU:CD1	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:CYS:SG	1:B:333:LEU:HD23	2.59	0.43
1:B:246:HIS:NE2	1:B:279:GLU:OE2	2.51	0.43
1:B:262:PRO:HA	1:B:263:PRO:HD3	1.83	0.43
1:B:15:GLU:O	1:B:16:LEU:C	2.61	0.42
1:A:122:GLU:OE1	1:A:159:PRO:HA	2.20	0.42
1:A:359:VAL:HG12	1:A:360:SER:N	2.34	0.42
1:A:276:SER:O	1:A:277:GLU:C	2.61	0.42
1:A:206:THR:OG1	1:A:234:VAL:HG22	2.20	0.42
1:A:9:THR:HG23	1:A:12:GLN:OE1	2.19	0.42
1:A:177:ILE:O	1:A:178:CYS:C	2.60	0.41
1:A:41:ALA:HB2	1:A:51:ASN:HD22	1.85	0.41
1:B:183:ILE:HD13	1:B:183:ILE:HG21	1.78	0.41
1:B:235:THR:HG22	1:B:249:ILE:HD13	2.02	0.41
1:B:78:ILE:HG12	1:B:145:PHE:CE1	2.55	0.41
1:B:109:VAL:HG21	1:B:174:TYR:CE1	2.52	0.41
1:A:52:THR:O	1:A:56:ARG:HG3	2.21	0.41
1:A:113:VAL:HA	1:A:124:THR:O	2.20	0.41
1:B:169:ASN:HD22	3:B:404:GOL:C3	2.34	0.41
1:A:122:GLU:HG2	4:A:531:HOH:O	2.19	0.41
1:B:285:ASN:O	1:B:286:ALA:C	2.64	0.41
1:B:309:SER:O	1:B:310:ALA:C	2.62	0.41
1:B:230:LYS:HD2	1:B:301:SER:CB	2.51	0.41
1:B:200:LYS:NZ	2:B:402:PO4:P	2.94	0.40
1:B:94:PHE:O	1:B:95:PRO:C	2.64	0.40
1:A:148:TRP:CD1	1:A:148:TRP:C	3.00	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	349/364 (96%)	332 (95%)	17 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	356/364 (98%)	340 (96%)	15 (4%)	1 (0%)	36	26
All	All	705/728 (97%)	672 (95%)	32 (4%)	1 (0%)	48	37

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	189	PRO

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	283/292 (97%)	282 (100%)	1 (0%)	84	79
1	B	286/292 (98%)	284 (99%)	2 (1%)	76	69
All	All	569/584 (97%)	566 (100%)	3 (0%)	81	76

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	296	TRP
1	B	11	GLU
1	B	296	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	307	GLN
1	B	45	GLN
1	B	96	GLN
1	B	181	ASN
1	B	283	ASN
1	B	320	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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
3	GOL	B	404	-	5,5,5	1.88	1 (20%)	5,5,5	1.96	2 (40%)
3	GOL	A	403	-	5,5,5	1.56	0	5,5,5	1.24	0
3	GOL	B	403	-	5,5,5	2.43	2 (40%)	5,5,5	1.08	0
2	PO4	B	401	-	4,4,4	1.53	1 (25%)	6,6,6	1.08	0
3	GOL	A	405	-	5,5,5	1.96	2 (40%)	5,5,5	1.60	2 (40%)
3	GOL	A	404	-	5,5,5	0.96	0	5,5,5	1.30	0
2	PO4	B	402	-	4,4,4	1.23	1 (25%)	6,6,6	0.77	0
3	GOL	B	405	-	5,5,5	2.34	3 (60%)	5,5,5	0.74	0
2	PO4	A	401	-	4,4,4	1.07	0	6,6,6	1.32	0
2	PO4	A	402	-	4,4,4	1.01	0	6,6,6	0.77	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	B	404	-	-	2/4/4/4	-
3	GOL	A	403	-	-	0/4/4/4	-
3	GOL	A	405	-	-	4/4/4/4	-
3	GOL	A	404	-	-	1/4/4/4	-
3	GOL	B	405	-	-	4/4/4/4	-
3	GOL	B	403	-	-	3/4/4/4	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	405	GOL	C3-C2	4.19	1.67	1.51
3	B	403	GOL	C3-C2	4.07	1.67	1.51
3	B	404	GOL	O1-C1	3.83	1.58	1.42
3	A	405	GOL	C3-C2	3.36	1.64	1.51
3	B	403	GOL	C1-C2	3.18	1.63	1.51
3	A	405	GOL	C1-C2	2.63	1.61	1.51
2	B	401	PO4	P-O1	2.54	1.56	1.50
2	B	402	PO4	P-O1	2.40	1.56	1.50
3	B	405	GOL	C1-C2	2.30	1.60	1.51
3	B	405	GOL	O2-C2	2.11	1.49	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	404	GOL	C3-C2-C1	-3.04	100.66	111.80
3	B	404	GOL	O1-C1-C2	-2.55	98.91	110.38
3	A	405	GOL	C3-C2-C1	-2.54	102.50	111.80
3	A	405	GOL	O2-C2-C3	2.04	117.62	109.18

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	405	GOL	C1-C2-C3-O3
3	B	404	GOL	C1-C2-C3-O3
3	B	405	GOL	O1-C1-C2-O2
3	B	405	GOL	O1-C1-C2-C3
3	A	405	GOL	O1-C1-C2-C3
3	B	403	GOL	C1-C2-C3-O3
3	B	405	GOL	C1-C2-C3-O3
3	A	405	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	B	403	GOL	O2-C2-C3-O3
3	B	404	GOL	O2-C2-C3-O3
3	B	405	GOL	O2-C2-C3-O3
3	A	404	GOL	O2-C2-C3-O3
3	B	403	GOL	O1-C1-C2-O2
3	A	405	GOL	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	404	GOL	3	0
3	B	403	GOL	1	0
3	A	405	GOL	1	0
3	A	404	GOL	2	0
2	B	402	PO4	2	0
3	B	405	GOL	3	0
2	A	401	PO4	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	353/364 (96%)	0.49	15 (4%) 40 44	27, 45, 64, 74	0
1	B	358/364 (98%)	0.53	20 (5%) 30 32	27, 45, 62, 90	0
All	All	711/728 (97%)	0.51	35 (4%) 35 37	27, 45, 64, 90	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	362	HIS	6.0
1	A	349	ALA	3.8
1	A	361	ASN	3.7
1	B	347	GLY	3.7
1	B	5	TYR	3.4
1	B	349	ALA	3.4
1	B	311	LEU	3.4
1	B	156	GLU	3.3
1	B	346	SER	3.3
1	B	348	GLN	3.1
1	A	156	GLU	3.0
1	B	7	ALA	3.0
1	B	326	GLU	2.8
1	B	157	HIS	2.8
1	A	6	PRO	2.8
1	A	345	PRO	2.7
1	A	39	SER	2.7
1	A	251	MET	2.6
1	B	304	ARG	2.5
1	A	291	PRO	2.5
1	B	98	CYS	2.4
1	A	62	LEU	2.4
1	A	43	ARG	2.4
1	A	342	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	344	THR	2.3
1	A	359	VAL	2.3
1	B	350	GLY	2.3
1	B	279	GLU	2.2
1	A	344	THR	2.2
1	B	174	TYR	2.0
1	A	96	GLN	2.0
1	B	6	PRO	2.0
1	B	293	LEU	2.0
1	A	318	LYS	2.0
1	B	93	PRO	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	405	6/6	0.60	0.21	61,74,83,93	0
3	GOL	A	403	6/6	0.75	0.17	48,64,78,85	0
3	GOL	B	403	6/6	0.76	0.20	47,61,76,80	0
3	GOL	A	405	6/6	0.80	0.19	43,52,57,60	14
2	PO4	A	401	5/5	0.81	0.15	47,50,60,66	0
2	PO4	B	401	5/5	0.83	0.13	51,53,53,65	0
3	GOL	B	404	6/6	0.84	0.15	37,53,64,65	0
3	GOL	A	404	6/6	0.88	0.18	62,75,80,89	14
2	PO4	B	402	5/5	0.93	0.08	52,54,64,73	0
2	PO4	A	402	5/5	0.96	0.06	47,56,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.