



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:09 AM GMT

PDB ID : 1ZY4  
Title : Crystal Structure of eIF2alpha Protein Kinase GCN2: R794G Hyperactivating Mutant in Apo Form.  
Authors : Padyana, A.K.; Qiu, H.; Roll-Mecak, A.; Hinnebusch, A.G.; Burley, S.K.  
Deposited on : 2005-06-09  
Resolution : 1.95 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

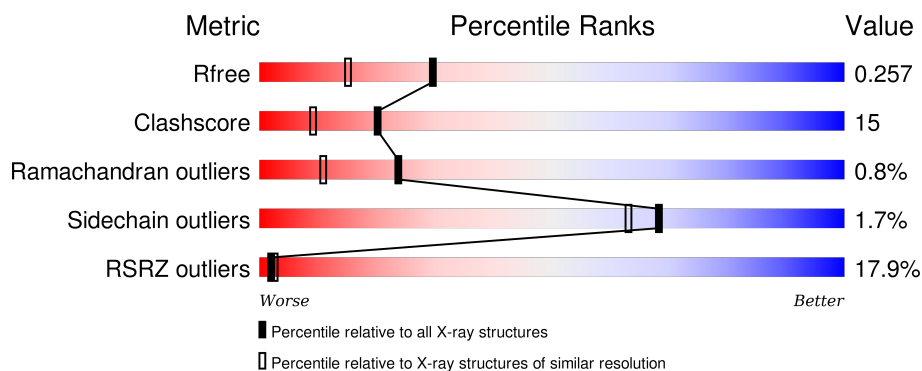
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	303	
1	B	303	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	998	-	X	-	-
2	GOL	B	998	-	X	-	-

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4539 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.

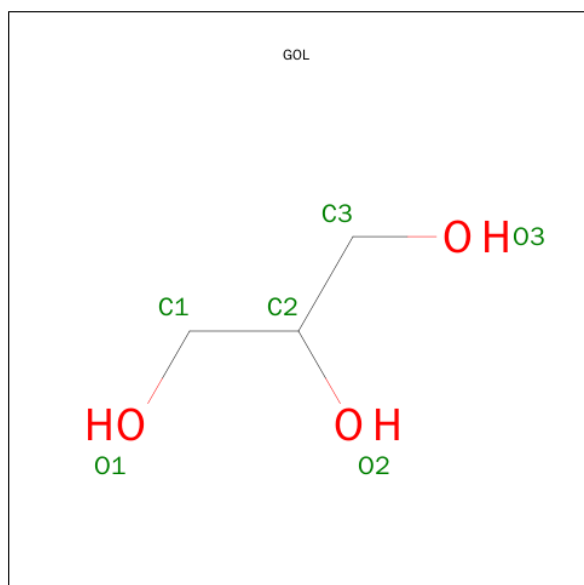
- Molecule 1 is a protein called Serine/threonine-protein kinase GCN2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2236	1434	390	403	9			
1	B	249	Total	C	N	O	S	0	0	0
			2057	1326	352	370	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	592	SER	-	CLONING ARTIFACT	UNP P15442
A	593	LEU	-	CLONING ARTIFACT	UNP P15442
A	794	GLY	ARG	ENGINEERED	UNP P15442
B	592	SER	-	CLONING ARTIFACT	UNP P15442
B	593	LEU	-	CLONING ARTIFACT	UNP P15442
B	794	GLY	ARG	ENGINEERED	UNP P15442

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	163	Total	O	0	0
			163	163		
3	B	71	Total	O	0	0
			71	71		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.83 Å 79.26 Å 146.86 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.72 – 1.95 37.72 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.9 (37.72-1.95) 99.1 (37.72-1.95)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.50 (at 1.95 Å)	Xtriage
Refinement program	CNX 2000.1	Depositor
R, $R_{free}$	0.224 , 0.252 0.222 , 0.257	Depositor DCC
$R_{free}$ test set	1407 reflections (3.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.2	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 57.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 45110 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4539	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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 [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	0.46	0/2281	0.65	0/3068
1	B	0.40	0/2098	0.58	0/2824
All	All	0.43	0/4379	0.62	0/5892

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	624	TYR	Sidechain

### 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	2236	0	2253	61	0
1	B	2057	0	2072	69	0
2	A	6	0	4	0	0
2	B	6	0	4	0	0
3	A	163	0	0	8	0
3	B	71	0	0	7	0
All	All	4539	0	4333	127	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 15.

All (127) 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:617:ARG:HH21	1:A:622:SER:HB3	1.17	1.04
1:B:618:ASN:HD21	1:B:620:LEU:HB2	1.35	0.90
1:A:657:ARG:HD3	3:A:95:HOH:O	1.74	0.86
1:A:953:VAL:HG12	3:A:102:HOH:O	1.79	0.81
1:B:958:ILE:O	1:B:962:ILE:HG12	1.81	0.81
1:B:652:HIS:HD2	1:B:654:TYR:H	1.28	0.79
1:B:833:HIS:HD2	1:B:835:ASP:H	1.29	0.79
1:B:790:TYR:CE2	1:B:792:GLU:HB3	2.19	0.78
1:A:617:ARG:NH2	1:A:622:SER:HB3	1.98	0.77
1:B:632:HIS:CE1	1:B:636:LYS:HD2	2.21	0.75
1:A:652:HIS:HD2	1:A:654:TYR:H	1.36	0.72
1:A:617:ARG:HH21	1:A:622:SER:CB	2.00	0.71
1:B:817:ARG:HH21	1:B:981:LEU:HB2	1.56	0.70
1:B:610:PHE:HB3	3:B:213:HOH:O	1.92	0.69
1:A:951:MET:HB3	1:A:954:GLU:HG2	1.75	0.69
1:B:595:TYR:OH	1:B:601:GLU:HB3	1.92	0.68
1:B:903:ASN:HD21	1:B:905:LYS:HB2	1.57	0.68
1:B:618:ASN:ND2	1:B:620:LEU:HB2	2.08	0.66
1:A:652:HIS:CD2	1:A:654:TYR:H	2.14	0.66
1:A:779:LYS:HG3	1:A:780:LYS:H	1.60	0.65
1:B:618:ASN:ND2	1:B:620:LEU:H	1.95	0.64
1:A:859:ASN:HD21	1:A:861:HIS:CD2	2.15	0.64
1:A:618:ASN:HD21	1:A:620:LEU:HB2	1.63	0.64
1:B:652:HIS:CD2	1:B:654:TYR:H	2.12	0.63
1:A:952:LYS:HE2	1:A:956:LYS:NZ	2.13	0.63
1:A:950:LYS:HB3	1:A:950:LYS:NZ	2.15	0.61
1:A:942:PHE:CE1	1:A:955:LYS:HG3	2.35	0.61
1:B:793:ASN:HB3	3:B:136:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:618:ASN:HD22	1:A:621:ASP:H	1.50	0.60
1:B:618:ASN:HD22	1:B:621:ASP:H	1.50	0.60
1:A:817:ARG:NH2	3:A:116:HOH:O	2.36	0.59
1:A:889:MET:HG2	1:A:932:LEU:CD1	2.32	0.59
1:A:951:MET:HB3	1:A:954:GLU:CG	2.33	0.57
1:B:804:ASN:O	1:B:807:GLN:HG2	2.05	0.56
1:A:845:GLU:HG2	3:A:64:HOH:O	2.04	0.56
1:B:942:PHE:CE1	1:B:955:LYS:HG3	2.41	0.56
1:B:793:ASN:HB2	3:B:87:HOH:O	2.06	0.56
1:B:931:ILE:HG23	3:B:195:HOH:O	2.05	0.56
1:A:618:ASN:ND2	1:A:620:LEU:H	2.03	0.56
1:A:960:LEU:O	1:A:969:ARG:HG3	2.05	0.56
1:B:615:LYS:HE2	1:B:624:TYR:CD2	2.41	0.55
1:B:833:HIS:HE1	1:B:852:GLY:O	1.91	0.54
1:A:645:MET:CE	1:B:784:LEU:HD11	2.38	0.54
1:A:889:MET:HG2	1:A:932:LEU:HD12	1.90	0.54
1:A:922:PHE:CD2	1:A:928:ARG:HA	2.43	0.53
1:A:942:PHE:CG	1:A:955:LYS:HE3	2.44	0.53
1:B:893:THR:HG21	1:B:964:HIS:CD2	2.44	0.52
1:B:910:SER:O	1:B:914:ILE:HG13	2.10	0.52
1:A:889:MET:HE3	1:A:889:MET:HA	1.91	0.52
1:A:893:THR:HG21	1:A:964:HIS:CE1	2.45	0.52
1:B:847:ARG:NH1	1:B:847:ARG:HB3	2.24	0.51
1:B:939:SER:HB3	1:B:941:GLU:HG3	1.92	0.51
1:B:954:GLU:O	1:B:958:ILE:HG13	2.10	0.51
1:B:618:ASN:HB3	1:B:621:ASP:OD2	2.10	0.51
1:A:780:LYS:HA	3:A:200:HOH:O	2.10	0.51
1:B:809:ARG:HA	1:B:812:TYR:CE2	2.46	0.51
1:B:904:GLU:N	1:B:904:GLU:OE2	2.39	0.51
1:A:861:HIS:N	1:A:861:HIS:ND1	2.59	0.50
1:B:611:GLY:O	1:B:612:GLN:HB3	2.12	0.50
1:B:656:VAL:HG13	1:B:789:GLU:HB3	1.92	0.50
1:B:817:ARG:HH21	1:B:981:LEU:CB	2.21	0.50
1:A:859:ASN:ND2	1:A:861:HIS:CD2	2.78	0.50
1:B:809:ARG:HH11	1:B:809:ARG:HG3	1.76	0.50
1:B:664:GLU:HB3	1:B:769:ARG:HH12	1.77	0.50
1:B:802:SER:O	1:B:803:GLU:HG3	2.12	0.50
1:A:952:LYS:HE2	1:A:956:LYS:HZ1	1.76	0.49
1:B:824:SER:OG	1:B:973:ARG:HD2	2.12	0.49
1:B:804:ASN:HB3	1:B:807:GLN:HG2	1.94	0.49
1:B:929:VAL:O	1:B:933:LYS:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:803:GLU:O	1:A:804:ASN:CG	2.51	0.49
1:A:634:GLU:HB3	3:A:217:HOH:O	2.13	0.49
1:A:771:PHE:CD1	1:A:780:LYS:HE2	2.47	0.49
1:B:813:TRP:HH2	1:B:957:ILE:HD12	1.78	0.48
1:A:950:LYS:HB3	1:A:950:LYS:HZ3	1.79	0.48
1:A:813:TRP:CE2	1:A:982:PRO:HG2	2.49	0.48
1:B:835:ASP:O	1:B:840:ASN:OD1	2.32	0.47
1:A:990:ILE:O	1:A:994:LEU:HG	2.13	0.47
1:B:834:ARG:HD2	1:B:855:GLY:O	2.15	0.47
1:B:973:ARG:HD3	3:B:180:HOH:O	2.14	0.47
1:B:595:TYR:HH	1:B:601:GLU:HB3	1.78	0.46
1:A:889:MET:HG2	1:A:932:LEU:HD13	1.97	0.46
1:B:946:PHE:O	1:B:948:ASP:N	2.48	0.46
1:A:607:GLN:HG3	1:A:612:GLN:HB3	1.97	0.46
1:B:889:MET:CE	1:B:936:ARG:HD3	2.45	0.46
1:B:885:ILE:O	1:B:885:ILE:HG23	2.15	0.46
1:A:641:LEU:O	1:A:645:MET:HG2	2.15	0.46
1:A:889:MET:HE2	1:A:936:ARG:HD3	1.98	0.45
1:B:638:SER:HA	1:B:641:LEU:HG	1.98	0.45
1:A:645:MET:HE1	1:B:784:LEU:HD11	1.98	0.45
1:B:940:ILE:HG22	1:B:940:ILE:O	2.16	0.45
1:A:889:MET:CE	1:A:936:ARG:HD3	2.47	0.45
1:A:845:GLU:CG	3:A:64:HOH:O	2.65	0.44
1:B:922:PHE:CD2	1:B:928:ARG:HA	2.52	0.44
1:B:903:ASN:H	1:B:903:ASN:HD22	1.65	0.44
1:A:803:GLU:O	1:A:804:ASN:C	2.54	0.44
1:B:977:ASN:HB3	3:B:156:HOH:O	2.17	0.44
1:A:961:LEU:O	1:A:969:ARG:HG2	2.18	0.44
1:A:833:HIS:O	1:A:834:ARG:HB2	2.17	0.44
1:B:847:ARG:HB3	1:B:847:ARG:HH11	1.82	0.44
1:A:771:PHE:CE1	1:A:780:LYS:HE2	2.53	0.43
1:B:931:ILE:O	1:B:934:LYS:HB2	2.18	0.43
1:A:897:ASP:OD1	1:A:898:GLY:N	2.51	0.43
1:B:661:ALA:HA	1:B:785:PHE:O	2.19	0.43
1:B:947:ASP:HB3	1:B:950:LYS:HB3	2.00	0.43
1:A:952:LYS:HE2	1:A:956:LYS:HZ3	1.80	0.43
1:B:804:ASN:HA	3:B:134:HOH:O	2.17	0.43
1:A:973:ARG:NH1	3:A:190:HOH:O	2.46	0.42
1:A:653:GLN:HG3	1:A:654:TYR:CD2	2.55	0.42
1:B:903:ASN:ND2	1:B:905:LYS:HB2	2.28	0.42
1:A:922:PHE:CE2	1:A:928:ARG:HA	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:645:MET:HE3	1:B:641:LEU:HD22	2.02	0.42
1:B:925:GLY:O	1:B:929:VAL:HG23	2.19	0.42
1:B:817:ARG:NH2	1:B:981:LEU:HB2	2.29	0.41
1:A:940:ILE:HD11	1:A:964:HIS:HD2	1.84	0.41
1:B:809:ARG:HA	1:B:812:TYR:CZ	2.55	0.41
1:B:959:ARG:HH21	1:B:959:ARG:HG3	1.84	0.41
1:A:956:LYS:HE3	1:A:956:LYS:HB2	1.77	0.41
1:A:958:ILE:O	1:A:962:ILE:HG12	2.19	0.41
1:A:799:LEU:HD12	1:A:799:LEU:HA	1.92	0.41
1:A:889:MET:HE3	1:A:936:ARG:CD	2.51	0.41
1:B:981:LEU:N	1:B:981:LEU:HD12	2.36	0.41
1:B:965:ASP:OD1	1:B:968:LYS:N	2.53	0.41
1:B:847:ARG:CB	1:B:847:ARG:HH11	2.34	0.40
1:A:956:LYS:HD2	1:A:980:TRP:CZ3	2.55	0.40
1:A:985:HIS:HD2	1:A:987:ASP:H	1.70	0.40
1:B:647:LEU:HD23	1:B:831:ILE:HD13	2.04	0.40
1:B:611:GLY:HA3	1:B:629:LYS:O	2.21	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	265/303 (88%)	254 (96%)	9 (3%)	2 (1%)	24	11
1	B	241/303 (80%)	222 (92%)	17 (7%)	2 (1%)	24	11
All	All	506/606 (84%)	476 (94%)	26 (5%)	4 (1%)	24	11

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	947	ASP

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Mol	Chain	Res	Type
1	A	899	THR
1	B	612	GLN
1	A	940	ILE

### 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	243/271 (90%)	239 (98%)	4 (2%)	70	66
1	B	224/271 (83%)	220 (98%)	4 (2%)	66	60
All	All	467/542 (86%)	459 (98%)	8 (2%)	68	63

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	859	ASN
1	A	861	HIS
1	A	889	MET
1	A	954	GLU
1	B	651	ASN
1	B	849	VAL
1	B	903	ASN
1	B	954	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	618	ASN
1	A	652	HIS
1	A	787	GLN
1	A	804	ASN
1	A	806	ASN
1	A	807	GLN
1	A	840	ASN

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Mol	Chain	Res	Type
1	A	859	ASN
1	A	930	ASN
1	A	977	ASN
1	B	618	ASN
1	B	651	ASN
1	B	652	HIS
1	B	787	GLN
1	B	793	ASN
1	B	829	GLN
1	B	833	HIS
1	B	840	ASN
1	B	903	ASN
1	B	964	HIS
1	B	977	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 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	GOL	A	998	-	5,5,5	4.79	5 (100%)	5,5,5	5.53	3 (60%)
2	GOL	B	998	-	5,5,5	4.76	5 (100%)	5,5,5	5.67	3 (60%)

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	A	998	-	-	0/4/4/4	0/0/0/0
2	GOL	B	998	-	-	0/4/4/4	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	998	GOL	C3-C2	-8.19	1.21	1.52
2	B	998	GOL	C3-C2	-7.99	1.21	1.52
2	A	998	GOL	C1-C2	-3.48	1.39	1.52
2	A	998	GOL	O2-C2	-3.35	1.33	1.43
2	B	998	GOL	C1-C2	-3.22	1.40	1.52
2	B	998	GOL	O2-C2	-2.88	1.34	1.43
2	A	998	GOL	O3-C3	2.94	1.55	1.42
2	B	998	GOL	O3-C3	3.39	1.57	1.42
2	A	998	GOL	O1-C1	3.98	1.59	1.42
2	B	998	GOL	O1-C1	4.41	1.61	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	998	GOL	O1-C1-C2	2.76	123.58	110.18
2	B	998	GOL	O1-C1-C2	3.22	125.81	110.18
2	B	998	GOL	O2-C2-C3	6.67	139.22	108.65
2	A	998	GOL	O2-C2-C3	6.73	139.53	108.65
2	A	998	GOL	O3-C3-C2	9.98	158.56	110.18
2	B	998	GOL	O3-C3-C2	10.26	159.96	110.18

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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	271/303 (89%)	0.83	29 (10%) 8 12	19, 37, 73, 91	0
1	B	249/303 (82%)	1.47	64 (25%) 1 1	23, 56, 87, 97	0
All	All	520/606 (85%)	1.13	93 (17%) 2 2	19, 47, 83, 97	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	896	LEU	13.2
1	A	899	THR	9.5
1	B	983	VAL	8.2
1	B	609	ALA	7.3
1	B	893	THR	6.6
1	B	610	PHE	5.9
1	B	981	LEU	5.9
1	B	885	ILE	5.7
1	B	895	VAL	5.7
1	B	603	ALA	5.3
1	B	902	TYR	5.3
1	B	938	VAL	5.1
1	A	900	GLY	5.0
1	B	809	ARG	5.0
1	B	980	TRP	4.9
1	A	901	HIS	4.8
1	A	926	MET	4.5
1	B	810	ASP	4.1
1	B	940	ILE	4.0
1	B	950	LYS	3.9
1	A	861	HIS	3.8
1	A	994	LEU	3.8
1	B	939	SER	3.8
1	B	942	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	845	GLU	3.7
1	B	887	THR	3.6
1	A	898	GLY	3.6
1	B	792	GLU	3.5
1	B	960	LEU	3.5
1	B	892	ALA	3.5
1	B	951	MET	3.3
1	B	602	ILE	3.3
1	B	805	LEU	3.3
1	A	990	ILE	3.3
1	A	896	LEU	3.3
1	B	937	SER	3.2
1	A	897	ASP	3.2
1	A	949	ASN	3.1
1	A	815	LEU	3.1
1	B	799	LEU	3.0
1	A	617	ARG	3.0
1	A	647	LEU	3.0
1	B	641	LEU	3.0
1	B	808	GLN	3.0
1	B	813	TRP	3.0
1	B	771	PHE	3.0
1	A	944	PRO	2.9
1	B	952	LYS	2.9
1	B	650	LEU	2.8
1	A	993	ALA	2.8
1	B	946	PHE	2.8
1	B	894	GLU	2.7
1	B	886	GLY	2.7
1	B	909	TYR	2.7
1	B	953	VAL	2.6
1	A	859	ASN	2.6
1	B	846	SER	2.6
1	B	802	SER	2.6
1	B	931	ILE	2.5
1	A	624	TYR	2.5
1	A	609	ALA	2.5
1	B	967	ASN	2.4
1	B	891	VAL	2.4
1	B	947	ASP	2.4
1	B	804	ASN	2.4
1	B	793	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	607	GLN	2.4
1	A	819	ILE	2.4
1	B	784	LEU	2.3
1	A	661	ALA	2.3
1	A	930	ASN	2.3
1	B	888	ALA	2.3
1	A	780	LYS	2.3
1	A	937	SER	2.2
1	B	943	PRO	2.2
1	B	982	PRO	2.2
1	B	807	GLN	2.2
1	B	949	ASN	2.2
1	B	964	HIS	2.2
1	B	971	GLY	2.2
1	A	610	PHE	2.2
1	A	646	LEU	2.2
1	A	799	LEU	2.1
1	B	957	ILE	2.1
1	B	847	ARG	2.1
1	B	604	VAL	2.1
1	B	648	ALA	2.1
1	A	924	THR	2.1
1	B	622	SER	2.0
1	A	849	VAL	2.0
1	B	966	PRO	2.0
1	B	979	GLY	2.0
1	B	962	ILE	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 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( $\text{\AA}^2$ )	Q<0.9
2	GOL	A	998	6/6	0.87	0.15	1.29	34,42,52,53	0
2	GOL	B	998	6/6	0.86	0.15	-0.01	62,67,70,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.