



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:04 PM UTC

PDB ID : 2OKK / pdb_00002okk
Title : The X-ray crystal structure of the 65kDa isoform of Glutamic Acid Decarboxylase (GAD65)
Authors : Buckle, A.M.; Fenalti, G.; Law, R.H.P.; Whisstock, J.C.
Deposited on : 2007-01-17
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

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)
Xtriage (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

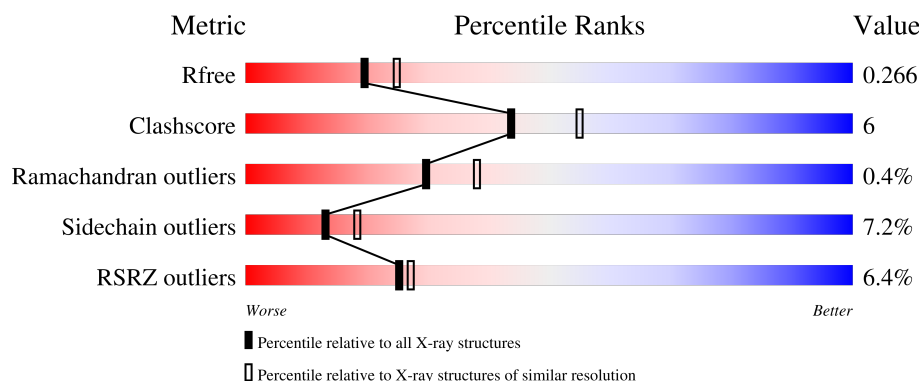
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 ≥ 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	497	

2 Entry composition [i](#)

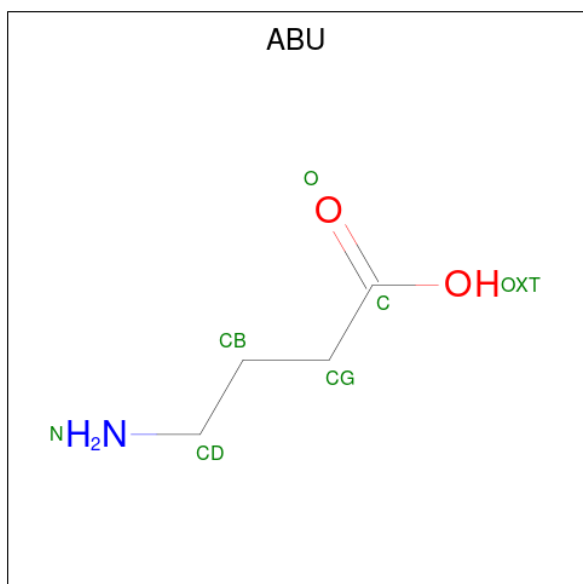
There are 4 unique types of molecules in this entry. The entry contains 3882 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutamate decarboxylase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	483	Total	C	N	O	P	S	0	0	0
			3770	2430	628	679	1	32			

- Molecule 2 is GAMMA-AMINO-BUTANOIC ACID (CCD ID: ABU) (formula: C₄H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			7	4	1	2		
2	A	1	Total	C	N	O	0	1
			7	4	1	2		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

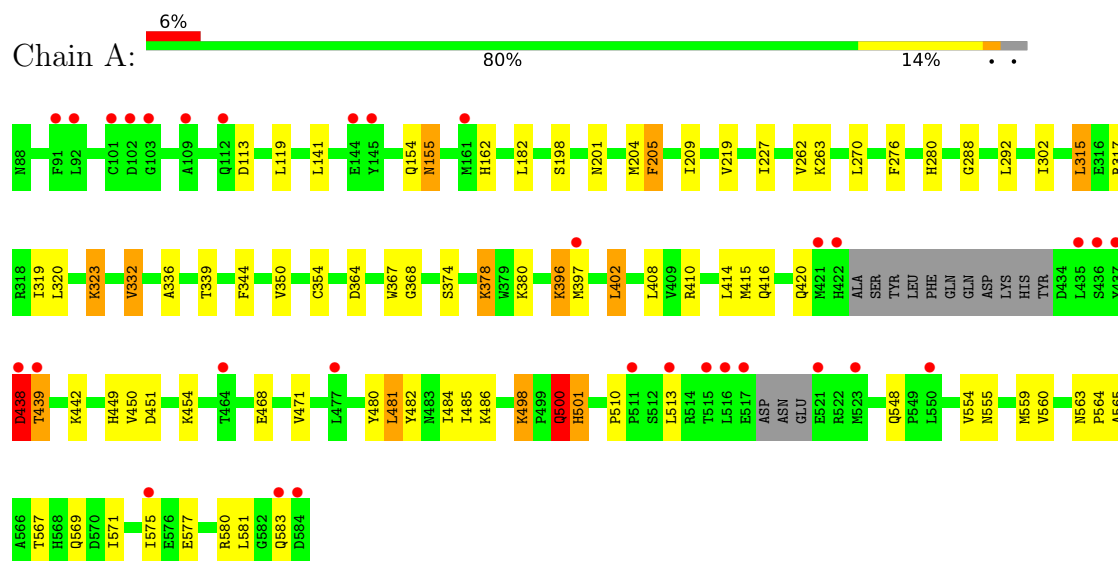
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	92	Total	O	0	0
			92	92		

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: Glutamate decarboxylase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	78.25Å 99.06Å 120.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.19 – 2.30 37.19 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (37.19-2.30) 97.9 (37.19-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.196 , 0.256 0.212 , 0.266	Depositor DCC
R_{free} test set	1500 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3882	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% of the height of the origin peak. No significant pseudotranslation is detected.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: LLP, ABU, 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	0.59	0/3834	0.95	10/5191 (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	A	1	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	500	GLN	N-CA-C	11.11	127.05	113.38
1	A	500	GLN	CA-C-N	7.45	135.11	121.70
1	A	500	GLN	C-N-CA	7.45	135.11	121.70
1	A	498	LYS	CA-C-N	6.41	126.43	119.89
1	A	498	LYS	C-N-CA	6.41	126.43	119.89
1	A	205	PHE	N-CA-C	6.19	118.83	111.71
1	A	438	ASP	CA-C-N	5.89	132.29	121.70
1	A	438	ASP	C-N-CA	5.89	132.29	121.70
1	A	402	LEU	N-CA-C	5.29	118.14	110.42
1	A	501	HIS	N-CA-C	5.25	125.70	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	501	HIS	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	500	GLN	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	3770	0	3697	46	0
2	A	14	0	0	0	0
3	A	6	0	8	0	0
4	A	92	0	0	3	0
All	All	3882	0	3705	46	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 (46) 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:438:ASP:HA	1:A:439:THR:OG1	1.52	1.08
1:A:481:LEU:HD11	1:A:559:MET:HG2	1.66	0.77
1:A:449:HIS:HE1	4:A:649:HOH:O	1.67	0.76
1:A:449:HIS:HD2	1:A:451:ASP:OD1	1.76	0.67
1:A:563:ASN:HD22	1:A:564:PRO:HD2	1.63	0.63
1:A:276:PHE:HB3	1:A:302:ILE:HD11	1.81	0.62
1:A:276:PHE:HB3	1:A:302:ILE:CD1	2.30	0.61
1:A:563:ASN:HD22	1:A:564:PRO:CD	2.16	0.58
1:A:336:ALA:O	1:A:368:GLY:HA3	2.04	0.58
1:A:481:LEU:CD1	1:A:559:MET:HG2	2.35	0.57
1:A:563:ASN:ND2	1:A:565:ALA:H	2.04	0.56
1:A:482:TYR:CZ	1:A:486:LYS:HD3	2.43	0.54
1:A:571:ILE:O	1:A:575:ILE:HG12	2.09	0.53
1:A:319:ILE:O	1:A:323:LYS:HG2	2.09	0.53
1:A:280:HIS:HD2	4:A:605:HOH:O	1.91	0.53
1:A:154:GLN:OE1	1:A:162:HIS:HE1	1.91	0.52
1:A:374:SER:HA	1:A:468:GLU:HG3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:396:LLP:O3	1:A:396:LLP:NZ	2.43	0.52
1:A:380:LYS:HD3	1:A:500:GLN:O	2.11	0.51
1:A:332:VAL:HG13	1:A:354:CYS:SG	2.50	0.50
1:A:364:ASP:OD2	1:A:396:LLP:N1	2.46	0.49
1:A:416:GLN:HE21	1:A:420:GLN:HE21	1.60	0.48
1:A:567:THR:OG1	1:A:569:GLN:HG2	2.13	0.48
1:A:367:TRP:CZ3	1:A:397:MET:HG2	2.49	0.47
1:A:155:ASN:HD22	1:A:155:ASN:C	2.23	0.46
1:A:480:TYR:CZ	1:A:484:ILE:HD11	2.50	0.46
1:A:482:TYR:O	1:A:486:LYS:HB2	2.16	0.46
1:A:262:VAL:HB	1:A:270:LEU:HD11	1.99	0.45
1:A:510:PRO:HD2	1:A:513:LEU:HD12	1.98	0.45
1:A:563:ASN:HD22	1:A:564:PRO:N	2.15	0.45
1:A:263:LYS:NZ	1:A:420:GLN:O	2.40	0.44
1:A:577:GLU:OE2	1:A:580:ARG:NH1	2.51	0.44
1:A:416:GLN:HE21	1:A:420:GLN:NE2	2.16	0.44
1:A:204:MET:SD	1:A:450:VAL:HG22	2.58	0.43
1:A:288:GLY:O	1:A:292:LEU:HG	2.18	0.43
1:A:315:LEU:HD22	1:A:319:ILE:HD11	2.00	0.43
1:A:454:LYS:NZ	4:A:664:HOH:O	2.21	0.43
1:A:438:ASP:CA	1:A:439:THR:OG1	2.44	0.42
1:A:415:MET:HB3	1:A:442:LYS:HG2	2.01	0.42
1:A:198:SER:O	1:A:201:ASN:HB2	2.20	0.42
1:A:344:PHE:H	1:A:500:GLN:NE2	2.19	0.41
1:A:378:LYS:HE3	1:A:378:LYS:HB2	1.89	0.41
1:A:397:MET:HE3	1:A:471:VAL:HG22	2.02	0.40
1:A:438:ASP:HA	1:A:439:THR:HG1	1.77	0.40
1:A:554:VAL:O	1:A:555:ASN:C	2.64	0.40
1:A:481:LEU:O	1:A:485:ILE:HG12	2.22	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	476/497 (96%)	462 (97%)	12 (2%)	2 (0%)	30	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	501	HIS
1	A	439	THR

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	387/424 (91%)	359 (93%)	28 (7%)	13	18

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	ASP
1	A	119	LEU
1	A	141	LEU
1	A	155	ASN
1	A	182	LEU
1	A	205	PHE
1	A	209	ILE
1	A	219	VAL
1	A	227	ILE
1	A	315	LEU
1	A	317	ARG
1	A	320	LEU
1	A	323	LYS
1	A	332	VAL
1	A	339	THR
1	A	350	VAL
1	A	378	LYS
1	A	402	LEU
1	A	408	LEU

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Mol	Chain	Res	Type
1	A	410	ARG
1	A	414	LEU
1	A	438	ASP
1	A	481	LEU
1	A	498	LYS
1	A	548	GLN
1	A	560	VAL
1	A	581	LEU
1	A	583	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	155	ASN
1	A	162	HIS
1	A	181	GLN
1	A	247	ASN
1	A	280	HIS
1	A	403	GLN
1	A	417	ASN
1	A	420	GLN
1	A	449	HIS
1	A	470	HIS
1	A	483	ASN
1	A	487	ASN
1	A	500	GLN
1	A	555	ASN
1	A	563	ASN
1	A	568	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	LLP	A	396	1	23,24,25	1.68	4 (17%)	25,32,34	1.79	4 (16%)

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	396	1	-	2/16/17/19	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	396	LLP	O3-C3	-5.57	1.24	1.36
1	A	396	LLP	C2-N1	2.66	1.38	1.33
1	A	396	LLP	C4-C4'	2.22	1.51	1.46
1	A	396	LLP	C4'-NZ	2.13	1.34	1.27

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	396	LLP	OP4-C5'-C5	5.26	119.21	109.36
1	A	396	LLP	OP4-P-OP1	-3.95	95.76	106.44
1	A	396	LLP	C4-C4'-NZ	-3.09	109.80	124.04
1	A	396	LLP	C5-C6-N1	-2.88	119.15	123.83

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	396	LLP	C4-C4'-NZ-CE
1	A	396	LLP	CD-CE-NZ-C4'

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	396	LLP	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	ABU	A	585[A]	-	6,6,6	0.87	0	6,6,6	0.82	0
3	GOL	A	587	-	5,5,5	0.37	0	5,5,5	0.16	0
2	ABU	A	586[B]	-	6,6,6	0.79	0	6,6,6	1.12	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	ABU	A	585[A]	-	-	1/4/4/4	-
3	GOL	A	587	-	-	2/4/4/4	-
2	ABU	A	586[B]	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	587	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	A	587	GOL	O1-C1-C2-O2
2	A	585[A]	ABU	CD-CB-CG-C
2	A	586[B]	ABU	CG-CB-CD-N
2	A	586[B]	ABU	OXT-C-CG-CB
2	A	586[B]	ABU	O-C-CG-CB

There are no ring outliers.

No monomer is involved in short contacts.

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	482/497 (96%)	0.81	31 (6%) 25 27	37, 51, 68, 96	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	521	GLU	5.4
1	A	91	PHE	4.9
1	A	584	ASP	4.6
1	A	437	TYR	4.0
1	A	421	MET	3.8
1	A	92	LEU	3.3
1	A	102	ASP	3.3
1	A	436	SER	3.2
1	A	439	THR	3.2
1	A	422	HIS	3.1
1	A	435	LEU	3.0
1	A	517	GLU	2.9
1	A	101	CYS	2.9
1	A	438	ASP	2.8
1	A	515	THR	2.6
1	A	550	LEU	2.5
1	A	523	MET	2.5
1	A	511	PRO	2.4
1	A	583	GLN	2.4
1	A	109	ALA	2.4
1	A	145	TYR	2.2
1	A	513	LEU	2.2
1	A	144	GLU	2.1
1	A	112	GLN	2.1
1	A	103	GLY	2.1
1	A	464	THR	2.1
1	A	397	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	575	ILE	2.1
1	A	161	MET	2.0
1	A	477	LEU	2.0
1	A	516	LEU	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. 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
1	LLP	A	396	24/25	0.96	0.08	41,44,52,55	0

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
2	ABU	A	586[B]	7/7	0.79	0.25	38,43,44,47	7
3	GOL	A	587	6/6	0.84	0.15	80,82,85,86	0
2	ABU	A	585[A]	7/7	0.87	0.19	32,35,36,36	7

6.5 Other polymers [i](#)

There are no such residues in this entry.