



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 4, 2026 – 10:37 PM UTC

PDB ID : 9NDR / pdb_00009ndr
Title : The 1.22 Angstrom crystal structure of galactose oxidase variant with genetically incorporated F2-Tyr495
Authors : Liu, A.; Li, J.; Graciano, A.
Deposited on : 2025-02-18
Resolution : 1.22 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

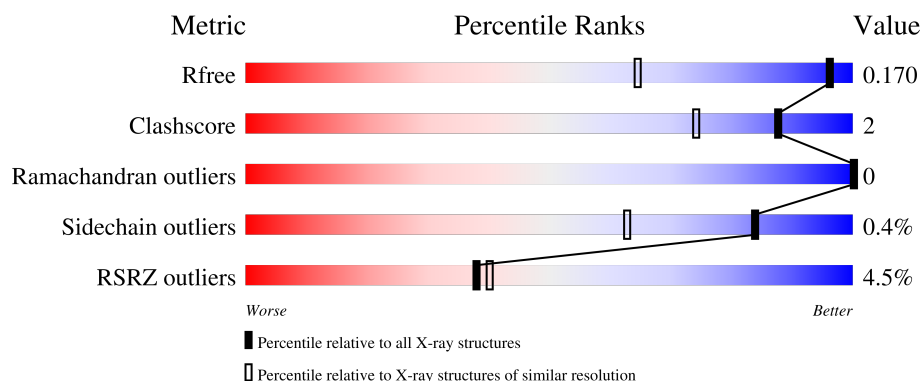
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.22 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	2002 (1.24-1.20)
Clashscore	190562	2061 (1.24-1.20)
Ramachandran outliers	187476	2009 (1.24-1.20)
Sidechain outliers	187428	2008 (1.24-1.20)
RSRZ outliers	180081	2000 (1.24-1.20)

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	648	<div> <div>4%</div> <div>94%</div> <div>5%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5706 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 Galactose oxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	640	Total	C	F	N	O	S	0	12	0
			4906	3066	2	853	966	19			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P0CS93
A	10	PRO	SER	conflict	UNP P0CS93
A	70	VAL	MET	conflict	UNP P0CS93
A	195	GLU	GLY	conflict	UNP P0CS93
A	494	ALA	VAL	conflict	UNP P0CS93
A	535	ASP	ASN	conflict	UNP P0CS93
A	640	LEU	-	expression tag	UNP P0CS93
A	641	GLU	-	expression tag	UNP P0CS93
A	642	HIS	-	expression tag	UNP P0CS93
A	643	HIS	-	expression tag	UNP P0CS93
A	644	HIS	-	expression tag	UNP P0CS93
A	645	HIS	-	expression tag	UNP P0CS93
A	646	HIS	-	expression tag	UNP P0CS93
A	647	HIS	-	expression tag	UNP P0CS93

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cu	0	0
			1	1		

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

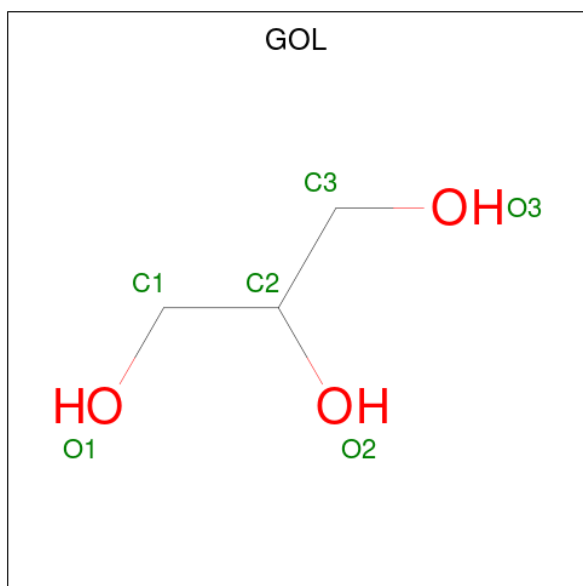
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

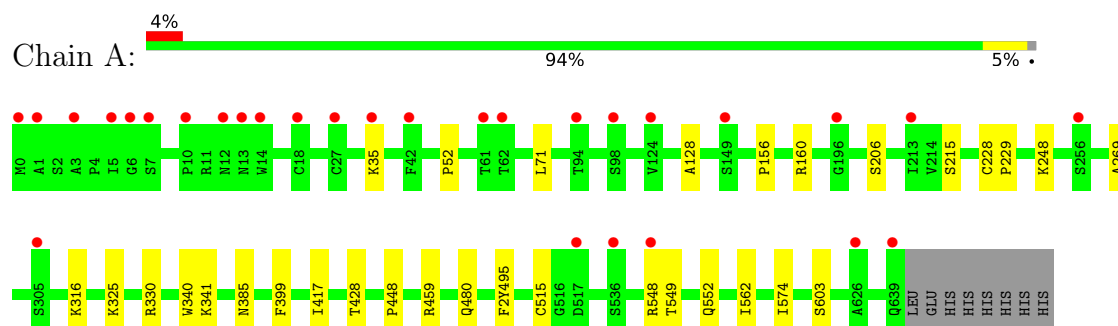
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	762	Total	O	0	0
			762	762		

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: Galactose oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.15Å 88.91Å 85.84Å 90.00° 117.97° 90.00°	Depositor
Resolution (Å)	48.28 – 1.22 48.28 – 1.22	Depositor EDS
% Data completeness (in resolution range)	98.0 (48.28-1.22) 95.2 (48.28-1.22)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 1.22Å)	Xtriage
Refinement program	PHENIX (1.21_5207-000)	Depositor
R, R_{free}	0.160 , 0.171 0.160 , 0.170	Depositor DCC
R_{free} test set	2000 reflections (1.06%)	wwPDB-VP
Wilson B-factor (Å ²)	14.4	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	5706	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CU, ACT, F2Y, CA, 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	0.29	2/5053 (0.0%)	0.52	2/6891 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	480[A]	GLN	CA-C	6.61	1.61	1.52
1	A	480[B]	GLN	CA-C	6.61	1.61	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	480[A]	GLN	CA-C-O	5.64	125.89	119.18
1	A	480[B]	GLN	CA-C-O	5.64	125.89	119.18

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4906	0	4701	19	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	12	0	9	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	24	0	32	2	0
6	A	762	0	0	5	0
All	All	5706	0	4742	19	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 2.

All (19) 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:325:LYS:NZ	6:A:801:HOH:O	2.21	0.73
1:A:160:ARG:HH21	5:A:706:GOL:H12	1.64	0.63
1:A:316:LYS:NZ	6:A:803:HOH:O	2.22	0.61
1:A:35:LYS:NZ	6:A:812:HOH:O	2.39	0.56
1:A:206:SER:OG	1:A:215:SER:OG	2.25	0.55
1:A:448:PRO:HA	1:A:574:ILE:HD11	1.90	0.53
1:A:459[B]:ARG:NH2	6:A:815:HOH:O	2.42	0.53
1:A:156:PRO:HA	5:A:706:GOL:H11	1.91	0.52
1:A:549:THR:HG22	1:A:562:ILE:HG22	1.93	0.51
1:A:228:CYS:N	1:A:229:PRO:HD3	2.28	0.48
1:A:548:ARG:HA	1:A:548:ARG:NE	2.29	0.47
1:A:35:LYS:HG2	1:A:71:LEU:HD21	1.97	0.47
1:A:417:ILE:HG12	1:A:428[B]:THR:HG22	1.97	0.46
1:A:340:TRP:CG	1:A:341:LYS:H	2.34	0.46
1:A:330:ARG:HH22	4:A:705:ACT:H1	1.82	0.45
1:A:248:LYS:HG3	1:A:269:ALA:HB2	2.00	0.43
1:A:385:ASN:HB2	1:A:399:PHE:CE2	2.55	0.42
1:A:603[B]:SER:OG	6:A:802:HOH:O	2.21	0.40
1:A:52:PRO:HA	1:A:128:ALA:O	2.21	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	649/648 (100%)	625 (96%)	24 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	538/534 (101%)	535 (99%)	3 (1%)	78	52

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

Mol	Chain	Res	Type
1	A	515	CYS
1	A	552[A]	GLN
1	A	552[B]	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	25	ASN
1	A	63	GLN
1	A	74	GLN
1	A	155	GLN
1	A	355	ASN
1	A	385	ASN
1	A	413	ASN
1	A	489	ASN
1	A	526	GLN

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	F2Y	A	495	1	13,14,15	1.03	1 (7%)	14,19,21	2.02	4 (28%)

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	F2Y	A	495	1	-	2/5/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	495	F2Y	OH-CZ	2.44	1.42	1.36

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	495	F2Y	CD1-CE1-CZ	-4.40	120.35	123.79
1	A	495	F2Y	F1-CE1-CZ	3.90	119.94	117.13
1	A	495	F2Y	CD2-CE2-CZ	-3.18	121.30	123.79
1	A	495	F2Y	F2-CE2-CZ	2.38	118.85	117.13

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	495	F2Y	CA-CB-CG-CD2
1	A	495	F2Y	CA-CB-CG-CD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 for Mogul analysis.

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	$\# Z > 2$
4	ACT	A	704	-	3,3,3	1.14	0	3,3,3	1.19	0
5	GOL	A	709	-	5,5,5	0.35	0	5,5,5	0.36	0
5	GOL	A	708	-	5,5,5	0.31	0	5,5,5	0.40	0
5	GOL	A	706	-	5,5,5	0.51	0	5,5,5	1.14	0
4	ACT	A	703	-	3,3,3	1.00	0	3,3,3	1.10	0
5	GOL	A	707	-	5,5,5	0.26	0	5,5,5	0.31	0
4	ACT	A	705	-	3,3,3	1.20	0	3,3,3	1.07	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
5	GOL	A	709	-	-	1/4/4/4	-
5	GOL	A	708	-	-	2/4/4/4	-
5	GOL	A	706	-	-	4/4/4/4	-
5	GOL	A	707	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	706	GOL	O1-C1-C2-C3
5	A	706	GOL	C1-C2-C3-O3
5	A	708	GOL	C1-C2-C3-O3
5	A	708	GOL	O2-C2-C3-O3
5	A	706	GOL	O1-C1-C2-O2
5	A	706	GOL	O2-C2-C3-O3
5	A	709	GOL	C1-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	706	GOL	2	0
4	A	705	ACT	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	639/648 (98%)	0.36	29 (4%) 38 40	7, 18, 33, 44	12 (1%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	MET	3.4
1	A	196	GLY	3.3
1	A	18	CYS	3.1
1	A	124	VAL	3.0
1	A	548	ARG	3.0
1	A	27	CYS	3.0
1	A	6	GLY	2.8
1	A	94	THR	2.8
1	A	12	ASN	2.7
1	A	626	ALA	2.7
1	A	14	TRP	2.7
1	A	639	GLN	2.7
1	A	517	ASP	2.5
1	A	536	SER	2.5
1	A	5	ILE	2.5
1	A	213	ILE	2.4
1	A	1	ALA	2.4
1	A	13	ASN	2.4
1	A	3	ALA	2.4
1	A	35	LYS	2.3
1	A	42	PHE	2.2
1	A	256	SER	2.2
1	A	305	SER	2.2
1	A	98	SER	2.2
1	A	10	PRO	2.2
1	A	61	THR	2.1
1	A	149	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	7	SER	2.0
1	A	62	THR	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	F2Y	A	495	14/15	0.97	0.06	13,15,18,18	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
4	ACT	A	704	4/4	0.61	0.18	34,35,36,38	0
4	ACT	A	705	4/4	0.66	0.19	31,32,32,33	0
5	GOL	A	708	6/6	0.75	0.16	42,45,45,46	0
5	GOL	A	707	6/6	0.83	0.13	24,26,27,30	0
5	GOL	A	706	6/6	0.84	0.14	21,28,34,34	0
5	GOL	A	709	6/6	0.91	0.11	22,27,31,34	0
3	CA	A	702	1/1	0.98	0.05	28,28,28,28	0
4	ACT	A	703	4/4	0.99	0.05	13,14,14,16	0
2	CU	A	701	1/1	1.00	0.02	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.