



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 16, 2020 – 01:13 PM EDT

PDB ID : 5MC2
Title : Crystal Structure of Gly278Asp mutant of Human Prolidase with Mn ions and GlyPro ligand
Authors : Wilk, P.; Mueller, U.; Dobbek, H.; Weiss, M.S.
Deposited on : 2016-11-09
Resolution : 1.70 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.10.1
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.10.1

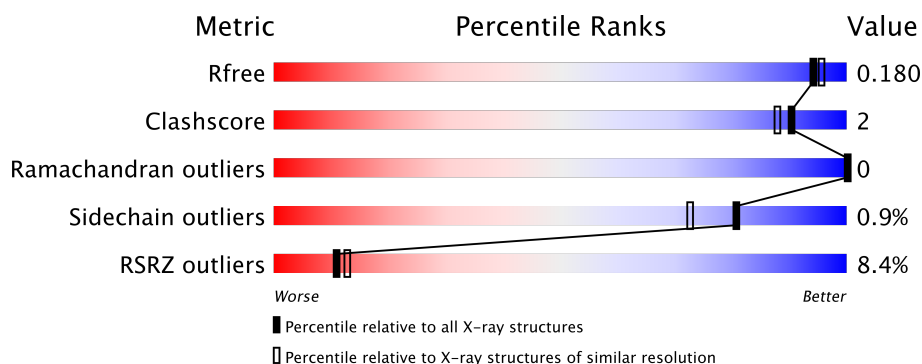
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	111664	3793 (1.70-1.70)
Clashscore	122126	4167 (1.70-1.70)
Ramachandran outliers	120053	4100 (1.70-1.70)
Sidechain outliers	120020	4100 (1.70-1.70)
RSRZ outliers	108989	3718 (1.70-1.70)

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 ≥ 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	484	<div> <div>9%</div> <div>94%</div> <div>5%</div> </div>
1	B	484	<div> <div>7%</div> <div>94%</div> <div>5%</div> </div>

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
6	GOL	A	505	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8757 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 Xaa-Pro dipeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	0	30	0
			3893	2473	674	715	31			
1	B	479	Total	C	N	O	S	0	30	0
			3905	2476	682	716	31			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	ALA	-	expression tag	UNP P12955
A	278	ASP	GLY	engineered mutation	UNP P12955
B	5	ALA	-	expression tag	UNP P12955
B	278	ASP	GLY	engineered mutation	UNP P12955

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

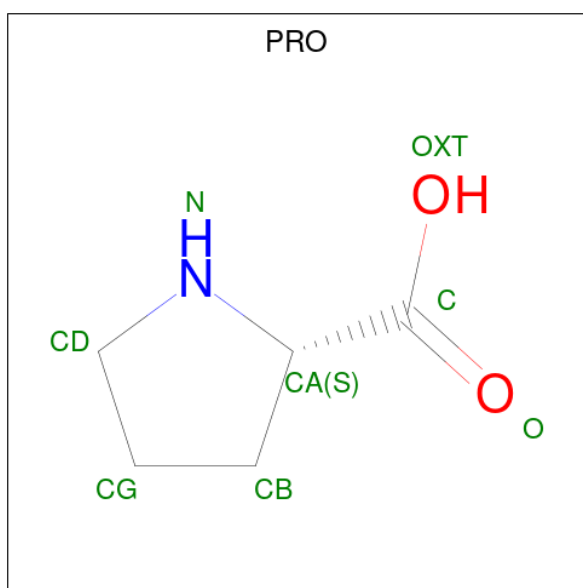
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is GLYCINE (three-letter code: GLY) (formula: C₂H₅NO₂).



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

- Molecule 5 is PROLINE (three-letter code: PRO) (formula: C₅H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	5	1	2		
5	B	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

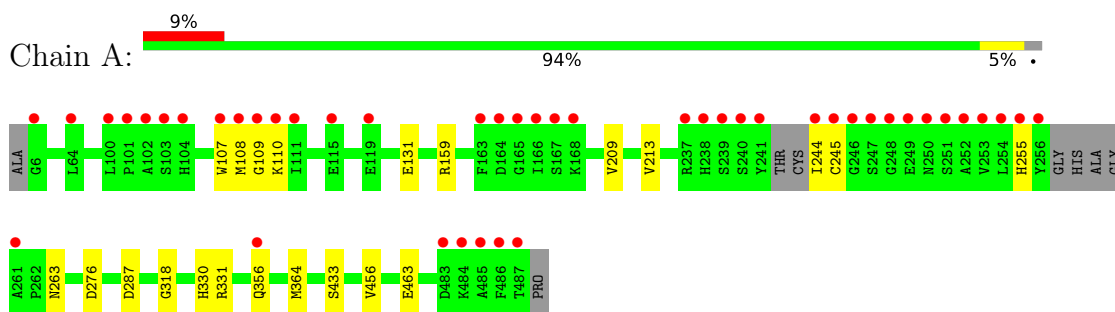
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	444	Total	O	0	4
			447	447		
7	B	433	Total	O	0	11
			442	442		

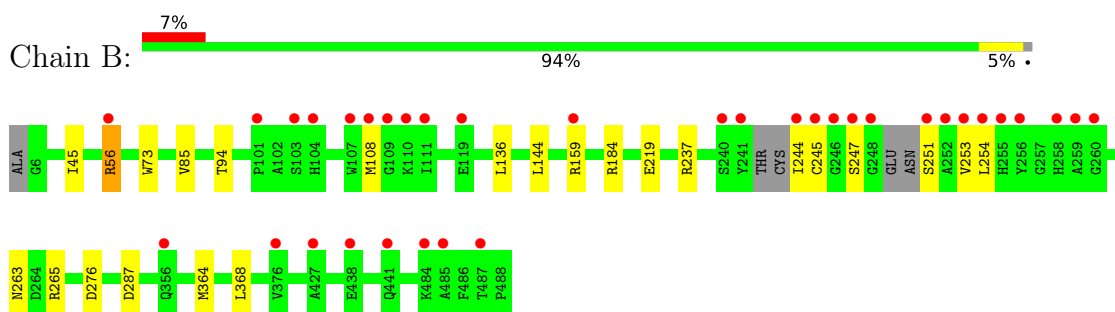
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Xaa-Pro dipeptidase



• Molecule 1: Xaa-Pro dipeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	103.84Å 108.25Å 211.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.17 – 1.70 48.17 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.17-1.70) 99.8 (48.17-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 1.70Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.141 , 0.171 0.151 , 0.180	Depositor DCC
R_{free} test set	1005 reflections (0.77%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.029 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8757	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: GOL, MN, NA

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	0.31	0/4066	0.51	0/5492
1	B	0.31	0/4077	0.51	0/5506
All	All	0.31	0/8143	0.51	0/10998

There are no bond length outliers.

There are no bond angle outliers.

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	3893	0	3877	17	0
1	B	3905	0	3896	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	4	0	2	0	0
4	B	4	0	2	0	0
5	A	8	0	7	0	0
5	B	8	0	7	0	0
6	A	18	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	24	0	32	0	0
7	A	447	0	0	7	0
7	B	442	0	0	5	0
All	All	8757	0	7847	30	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 (30) 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:B:287:ASP:OD1	7:B:601:HOH:O	2.10	0.70
1:B:276:ASP:OD1	7:B:601:HOH:O	2.12	0.68
1:B:247:SER:OG	1:B:265:ARG:NH2	2.28	0.65
1:A:287:ASP:OD1	7:A:601:HOH:O	2.15	0.63
1:A:131:GLU:OE2	7:A:602:HOH:O	2.15	0.63
1:A:331[B]:ARG:NH1	7:A:610:HOH:O	2.35	0.60
1:A:276:ASP:OD1	7:A:601:HOH:O	2.16	0.59
1:A:209:VAL:O	1:A:213[A]:VAL:HG23	2.06	0.55
1:A:456[B]:VAL:HG22	1:A:463:GLU:HB3	1.87	0.55
1:A:433[B]:SER:OG	7:A:603:HOH:O	2.20	0.47
1:B:244:ILE:HB	1:B:276:ASP:HB3	1.96	0.47
1:A:108:MET:HB3	1:B:253:VAL:HG11	1.96	0.47
1:A:244:ILE:HB	1:A:276:ASP:HB3	1.97	0.46
1:B:45[B]:ILE:HD11	1:B:144:LEU:HD22	1.99	0.45
1:B:159[B]:ARG:NH2	7:B:603:HOH:O	2.50	0.45
1:A:255:HIS:HB2	1:B:108:MET:CE	2.47	0.45
1:A:318:GLY:O	7:A:604:HOH:O	2.21	0.45
1:A:356[A]:GLN:NE2	7:A:614:HOH:O	2.49	0.45
1:B:219:GLU:OE1	1:B:251:SER:OG	2.28	0.45
1:B:85:VAL:HG22	1:B:94:THR:HB	1.98	0.44
1:B:364[A]:MET:HE3	1:B:368:LEU:HA	2.00	0.44
1:A:109:GLY:HA3	1:B:254:LEU:HD21	2.00	0.43
1:A:109:GLY:HA3	1:B:254:LEU:CD2	2.48	0.43
1:A:110:LYS:H	1:B:254:LEU:CD2	2.31	0.42
1:B:73:TRP:CD1	1:B:184:ARG:HG2	2.54	0.42
1:A:330:HIS:CE1	1:A:364[A]:MET:HE2	2.54	0.42
1:B:45[B]:ILE:CD1	1:B:136:LEU:HD22	2.50	0.42
1:B:56[A]:ARG:NH2	7:B:619:HOH:O	2.52	0.42
1:B:56[B]:ARG:NH2	7:B:613:HOH:O	2.51	0.41
1:A:107[A]:TRP:CE3	1:A:107[A]:TRP:HA	2.56	0.41

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	500/484 (103%)	492 (98%)	8 (2%)	0	100	100
1	B	503/484 (104%)	493 (98%)	10 (2%)	0	100	100
All	All	1003/968 (104%)	985 (98%)	18 (2%)	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	431/405 (106%)	428 (99%)	3 (1%)	85	79
1	B	431/405 (106%)	426 (99%)	5 (1%)	74	62
All	All	862/810 (106%)	854 (99%)	8 (1%)	81	73

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

Mol	Chain	Res	Type
1	A	159	ARG
1	A	245	CYS
1	A	263	ASN
1	B	56[A]	ARG
1	B	56[B]	ARG

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Mol	Chain	Res	Type
1	B	237	ARG
1	B	245	CYS
1	B	263	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	$\# Z > 2$
6	GOL	B	506	-	5,5,5	0.37	0	5,5,5	0.22	0
6	GOL	B	508	-	5,5,5	0.37	0	5,5,5	0.21	0
6	GOL	B	505	-	5,5,5	0.34	0	5,5,5	0.34	0
6	GOL	A	505	-	5,5,5	0.42	0	5,5,5	0.15	0
6	GOL	A	506	-	5,5,5	0.36	0	5,5,5	0.34	0
6	GOL	B	507	-	5,5,5	0.43	0	5,5,5	0.34	0
6	GOL	A	507	-	5,5,5	0.32	0	5,5,5	0.28	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
6	GOL	B	506	-	-	4/4/4/4	-
6	GOL	B	508	-	-	2/4/4/4	-
6	GOL	B	505	-	-	0/4/4/4	-
6	GOL	A	505	-	-	4/4/4/4	-
6	GOL	A	506	-	-	1/4/4/4	-
6	GOL	B	507	-	-	4/4/4/4	-
6	GOL	A	507	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	506	GOL	O1-C1-C2-C3
6	A	505	GOL	O1-C1-C2-C3
6	B	508	GOL	O1-C1-C2-C3
6	B	507	GOL	C1-C2-C3-O3
6	B	506	GOL	O1-C1-C2-O2
6	B	507	GOL	O2-C2-C3-O3
6	B	506	GOL	C1-C2-C3-O3
6	A	505	GOL	C1-C2-C3-O3
6	B	507	GOL	O1-C1-C2-C3
6	A	505	GOL	O2-C2-C3-O3
6	B	508	GOL	O1-C1-C2-O2
6	A	505	GOL	O1-C1-C2-O2
6	A	506	GOL	O1-C1-C2-C3
6	B	506	GOL	O2-C2-C3-O3
6	B	507	GOL	O1-C1-C2-O2

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, 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	476/484 (98%)	0.36	45 (9%) 8 9	20, 28, 77, 107	0
1	B	479/484 (98%)	0.09	35 (7%) 15 17	18, 27, 75, 102	0
All	All	955/968 (98%)	0.22	80 (8%) 11 13	18, 28, 77, 107	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	107[A]	TRP	29.2
1	B	244	ILE	19.9
1	A	485	ALA	13.0
1	B	241	TYR	12.2
1	A	244	ILE	8.0
1	A	247	SER	7.1
1	A	237[A]	ARG	7.1
1	A	101	PRO	7.0
1	A	249	GLU	6.4
1	A	486	PHE	6.4
1	A	108	MET	6.3
1	A	250	ASN	6.2
1	A	256	TYR	5.9
1	A	241	TYR	5.8
1	A	252	ALA	5.7
1	A	261	ALA	5.7
1	A	484	LYS	5.0
1	A	109	GLY	4.9
1	A	487	THR	4.8
1	B	247	SER	4.7
1	B	104	HIS	4.6
1	B	254	LEU	4.5
1	B	487	THR	4.5
1	A	254	LEU	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	251	SER	4.3
1	B	253	VAL	4.1
1	B	245	CYS	4.0
1	B	246	GLY	4.0
1	A	102	ALA	4.0
1	B	108	MET	3.8
1	B	255	HIS	3.7
1	A	248	GLY	3.7
1	B	111	ILE	3.7
1	A	245	CYS	3.6
1	B	252	ALA	3.6
1	B	107	TRP	3.6
1	B	485	ALA	3.6
1	B	240	SER	3.6
1	A	110	LYS	3.5
1	B	101	PRO	3.5
1	A	483	ASP	3.4
1	A	246	GLY	3.3
1	A	253	VAL	3.3
1	A	167	SER	3.2
1	A	115[A]	GLU	3.2
1	A	104	HIS	3.1
1	A	168	LYS	3.1
1	A	164[A]	ASP	3.0
1	B	259	ALA	3.0
1	B	56[A]	ARG	2.9
1	B	109	GLY	2.9
1	B	110	LYS	2.8
1	B	248	GLY	2.8
1	B	256	TYR	2.8
1	A	240	SER	2.7
1	A	165	GLY	2.6
1	A	6	GLY	2.6
1	B	159[A]	ARG	2.6
1	B	441	GLN	2.6
1	A	255	HIS	2.5
1	A	239	SER	2.5
1	B	103	SER	2.5
1	B	260	GLY	2.4
1	A	119	GLU	2.4
1	A	356[A]	GLN	2.4
1	B	119	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	64	LEU	2.4
1	A	103	SER	2.3
1	A	100	LEU	2.3
1	A	166	ILE	2.3
1	A	163	PHE	2.2
1	B	484	LYS	2.2
1	B	438	GLU	2.2
1	B	251	SER	2.2
1	B	427	ALA	2.1
1	A	238	HIS	2.1
1	B	356	GLN	2.1
1	A	111	ILE	2.1
1	B	376[A]	VAL	2.0
1	B	258	HIS	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. 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
6	GOL	B	505	6/6	0.43	0.39	82,83,84,84	0
6	GOL	A	506	6/6	0.67	0.33	82,83,83,83	0
6	GOL	A	505	6/6	0.70	0.42	83,84,84,85	0
6	GOL	B	508	6/6	0.71	0.18	52,58,59,60	0
6	GOL	B	507	6/6	0.71	0.28	74,77,77,78	0
6	GOL	A	507	6/6	0.75	0.16	43,49,50,51	0
6	GOL	B	506	6/6	0.87	0.32	74,74,74,74	0
4	GLY	B	503	4/5	0.96	0.18	26,28,30,33	0
5	PRO	B	504	8/8	0.97	0.10	26,26,29,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	PRO	A	504	8/8	0.97	0.08	25,28,31,32	0
3	NA	A	500	1/1	0.97	0.14	36,36,36,36	1
3	NA	B	500	1/1	0.98	0.22	33,33,33,33	1
4	GLY	A	503	4/5	0.98	0.13	27,28,31,32	0
2	MN	B	499	1/1	1.00	0.07	24,24,24,24	1
2	MN	A	499	1/1	1.00	0.09	23,23,23,23	1

6.5 Other polymers [i](#)

There are no such residues in this entry.