



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:16 PM UTC

PDB ID : 6SRF / pdb_00006srf
Title : Crystal Structure of Human Prolidase G278N variant expressed in the presence of chaperones
Authors : Wator, E.; Wilk, P.
Deposited on : 2019-09-05
Resolution : 1.85 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

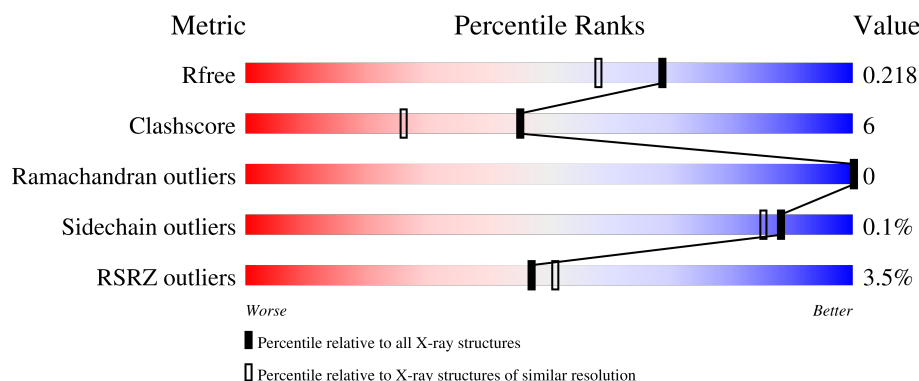
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

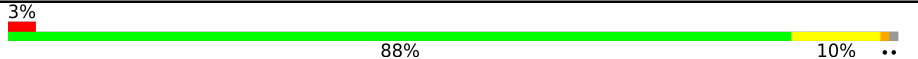
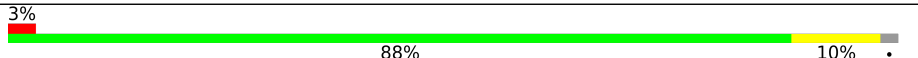
The reported resolution of this entry is 1.85 Å.

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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	488	
1	B	488	

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
4	GOL	A	503	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 16219 atoms, of which 7509 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	482	Total	C	H	N	O	S	0	12	0
			7542	2416	3721	663	710	32			
1	B	479	Total	C	H	N	O	S	0	19	0
			7577	2416	3756	660	713	32			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	278	ASP	GLY	engineered mutation	UNP P12955
B	278	ASP	GLY	engineered mutation	UNP P12955

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

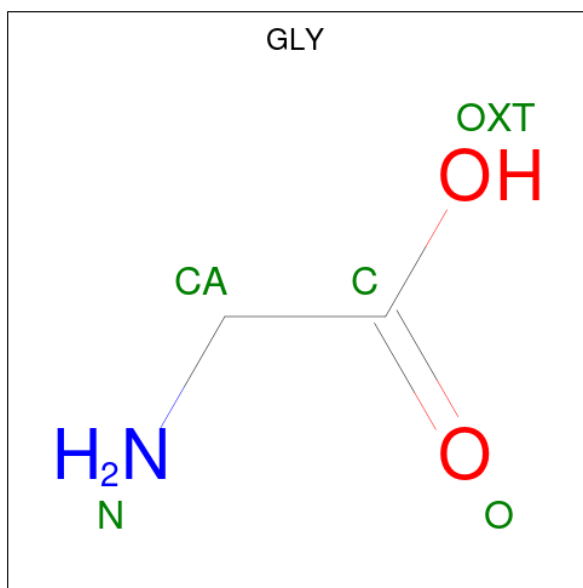
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



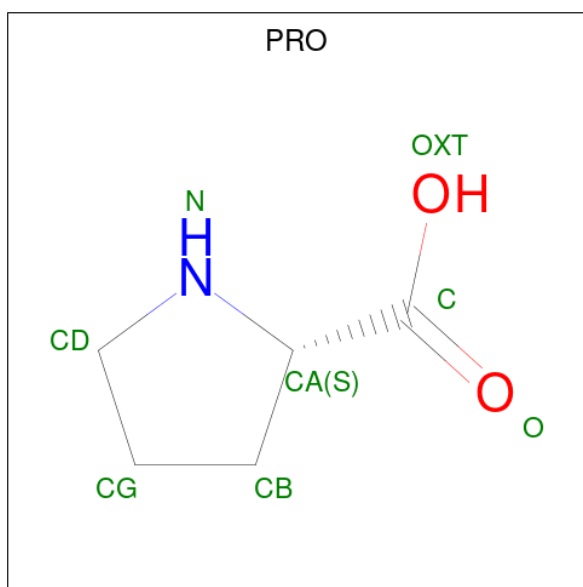
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	0	0
			9	2	5	1	1		
5	B	1	Total	C	H	N	O	0	0
			9	2	5	1	1		

- Molecule 6 is PROLINE (CCD ID: PRO) (formula: $C_5H_9NO_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	H	N	O	0	0
			15	5	7	1	2		
6	B	1	Total	C	H	N	O	0	0
			15	5	7	1	2		

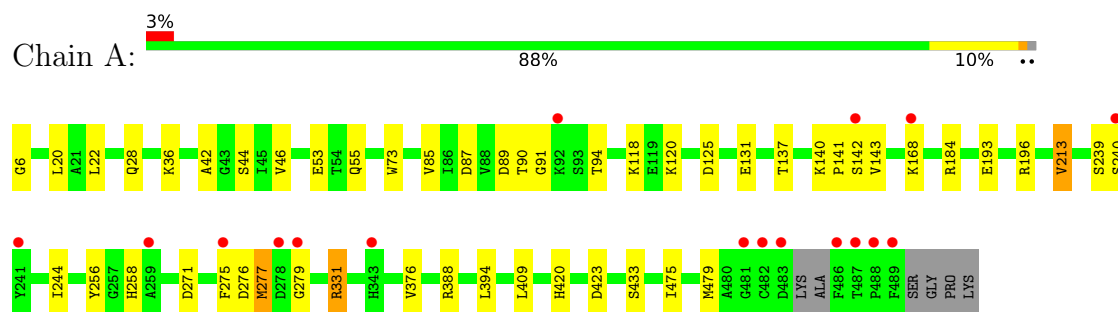
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	471	Total	O	0	0
			471	471		
7	B	563	Total	O	0	0
			563	563		

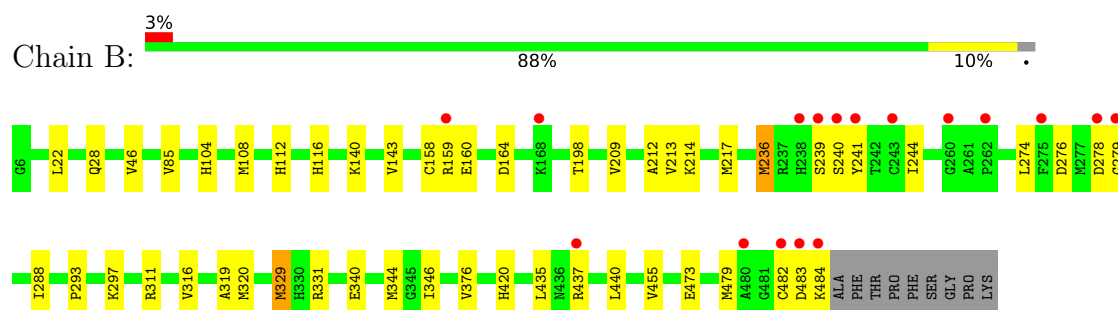
3 Residue-property plots [i](#)

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: 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	103.56Å 106.69Å 216.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.86 – 1.85 47.86 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.86-1.85) 99.6 (47.86-1.85)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 1.84Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.180 , 0.216 0.181 , 0.218	Depositor DCC
R_{free} test set	2101 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.407	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 56.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16219	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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: NA, MN, 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	0.39	4/3950 (0.1%)	0.49	5/5341 (0.1%)
1	B	0.59	9/3976 (0.2%)	0.54	4/5377 (0.1%)
All	All	0.50	13/7926 (0.2%)	0.52	9/10718 (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	2
1	B	0	1
All	All	0	3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	160	GLU	C-O	-14.51	1.06	1.23
1	B	159	ARG	C-O	-12.07	1.09	1.24
1	B	158	CYS	C-O	-10.12	1.11	1.23
1	B	160	GLU	CD-OE1	-7.87	1.10	1.25
1	B	160	GLU	CD-OE2	-7.55	1.11	1.25
1	A	331[A]	ARG	C-O	7.43	1.32	1.24
1	A	331[B]	ARG	C-O	7.43	1.32	1.24
1	B	329[A]	MET	N-CA	6.79	1.54	1.46
1	B	329[B]	MET	N-CA	6.79	1.54	1.46
1	B	297	LYS	C-O	-5.63	1.17	1.23
1	A	213	VAL	CB-CG2	5.48	1.70	1.52
1	A	213	VAL	CA-C	5.35	1.58	1.52
1	B	159	ARG	C-N	-5.10	1.26	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	437	ARG	CA-C-O	8.28	132.34	120.51
1	B	329[A]	MET	CA-C-O	6.47	127.41	119.97
1	B	329[B]	MET	CA-C-O	6.47	127.41	119.97
1	B	437	ARG	O-C-N	-5.39	115.42	122.59
1	A	331[A]	ARG	N-CA-C	5.33	117.09	111.28
1	A	331[B]	ARG	N-CA-C	5.33	117.09	111.28
1	A	331[A]	ARG	CA-C-O	5.27	126.06	120.10
1	A	331[B]	ARG	CA-C-O	5.27	126.06	120.10
1	A	409	LEU	CA-CB-CG	-5.03	98.71	116.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	213	VAL	Mainchain
1	A	277[B]	MET	Mainchain
1	B	236	MET	Mainchain

5.2 Too-close contacts ⓘ

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	3821	3721	3741	46	0
1	B	3821	3756	3757	41	1
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	6	8	8	5	0
5	A	4	5	2	0	0
5	B	4	5	2	0	0
6	A	8	7	7	0	0
6	B	8	7	7	0	0
7	A	471	0	0	23	0
7	B	563	0	0	10	0
All	All	8710	7509	7524	89	1

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 (89) 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:484:LYS:HA	7:B:967:HOH:O	1.61	0.99
1:A:168:LYS:NZ	7:A:603:HOH:O	1.94	0.98
1:A:137:THR:OG1	7:A:602:HOH:O	1.82	0.96
1:A:131:GLU:OE2	7:A:601:HOH:O	1.82	0.95
1:B:420:HIS:HD2	7:B:725:HOH:O	1.56	0.88
1:A:140:LYS:HA	1:A:140:LYS:HE2	1.58	0.85
4:A:503:GOL:H32	7:A:611:HOH:O	1.79	0.82
1:B:316:VAL:HA	1:B:329[A]:MET:CE	2.10	0.82
1:A:420:HIS:HD2	7:A:728:HOH:O	1.64	0.81
1:B:473:GLU:OE2	7:B:601:HOH:O	1.97	0.80
4:A:503:GOL:H32	7:A:742:HOH:O	1.83	0.79
1:A:55[A]:GLN:OE1	7:A:604:HOH:O	2.04	0.75
1:B:316:VAL:HA	1:B:329[A]:MET:HE1	1.73	0.71
1:A:140:LYS:NZ	7:A:609:HOH:O	2.18	0.69
4:A:503:GOL:C3	7:A:611:HOH:O	2.39	0.69
1:A:276:ASP:OD1	7:A:605:HOH:O	2.11	0.68
1:B:420:HIS:CD2	7:B:725:HOH:O	2.37	0.68
1:B:331[B]:ARG:NH2	7:B:605:HOH:O	2.26	0.68
1:B:316:VAL:HG13	1:B:329[A]:MET:CE	2.24	0.67
1:A:271:ASP:OD1	7:A:606:HOH:O	2.12	0.67
1:A:20:LEU:HD13	1:A:120:LYS:O	1.94	0.66
1:A:89:ASP:OD2	7:A:607:HOH:O	2.14	0.65
1:B:316:VAL:HG13	1:B:329[A]:MET:HE2	1.79	0.64
1:A:388:ARG:HB2	4:A:503:GOL:H31	1.81	0.63
1:A:331[A]:ARG:NH2	7:A:613:HOH:O	2.30	0.63
1:A:168:LYS:HD2	7:A:602:HOH:O	1.99	0.62
1:A:239:SER:HA	1:A:279:GLY:HA2	1.82	0.62
1:B:164[B]:ASP:OD2	7:B:603:HOH:O	2.16	0.60
1:A:420:HIS:CD2	7:A:728:HOH:O	2.44	0.60
1:A:423:ASP:OD2	7:A:608:HOH:O	2.16	0.58
1:B:239:SER:HA	1:B:279:GLY:HA2	1.86	0.58
1:B:241:TYR:OH	7:B:602:HOH:O	2.15	0.57
1:A:42:ALA:O	1:A:142:SER:HB3	2.06	0.55
1:A:20:LEU:CD1	1:A:120:LYS:O	2.54	0.55
1:A:433[B]:SER:OG	7:A:610:HOH:O	2.18	0.55
1:B:319:ALA:HB3	1:B:329[A]:MET:HE1	1.87	0.55
1:A:240:SER:HB3	1:A:376[B]:VAL:CG2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:483:ASP:O	1:B:484:LYS:HG3	2.09	0.52
1:B:316:VAL:HG13	1:B:329[A]:MET:HE3	1.92	0.51
1:B:28:GLN:NE2	7:B:604:HOH:O	2.25	0.51
1:A:140:LYS:HE2	1:A:140:LYS:CA	2.37	0.51
1:B:344:MET:HE2	1:B:346:ILE:HD12	1.91	0.51
1:A:275:PHE:HB3	1:A:277[A]:MET:HE2	1.93	0.51
1:A:87:ASP:O	1:A:91:GLY:N	2.40	0.50
1:B:244:ILE:HB	1:B:276:ASP:HB3	1.93	0.50
1:A:244:ILE:HB	1:A:276:ASP:HB3	1.92	0.50
1:A:193:GLU:OE1	1:A:196:ARG:NH1	2.42	0.50
1:B:278:ASP:CG	1:B:288:ILE:H	2.20	0.50
1:A:28:GLN:NE2	7:A:628:HOH:O	2.46	0.48
1:B:311:ARG:NH1	1:B:340:GLU:OE2	2.46	0.48
1:B:214:LYS:N	1:B:217:MET:HE3	2.26	0.48
1:A:331[B]:ARG:CZ	1:A:394:LEU:HD11	2.42	0.48
1:B:293:PRO:O	1:B:479:MET:HE1	2.14	0.48
1:A:53:GLU:CG	7:A:612:HOH:O	2.62	0.47
1:B:320:MET:HE1	1:B:455:VAL:HG21	1.97	0.47
1:B:316:VAL:CA	1:B:329[A]:MET:CE	2.88	0.46
1:B:241:TYR:OH	1:B:276:ASP:OD1	2.27	0.46
1:A:6:GLY:N	7:A:633:HOH:O	2.49	0.46
1:B:276:ASP:OD1	7:B:602:HOH:O	2.20	0.45
1:A:142:SER:OG	1:A:143:VAL:N	2.49	0.45
1:B:112:HIS:HB3	1:B:116:HIS:CD2	2.51	0.45
1:A:46:VAL:O	1:A:85:VAL:HA	2.17	0.45
1:A:118:LYS:NZ	7:A:630:HOH:O	2.48	0.45
1:B:240:SER:HB3	1:B:376[B]:VAL:CG2	2.47	0.44
1:A:87:ASP:OD2	1:A:90:THR:OG1	2.20	0.44
1:B:209:VAL:O	1:B:213[A]:VAL:HG23	2.18	0.44
1:B:22:LEU:C	1:B:22:LEU:HD23	2.43	0.44
1:B:214:LYS:H	1:B:217:MET:HE3	1.83	0.43
1:B:316:VAL:HG22	1:B:329[A]:MET:HE3	1.99	0.43
1:A:22:LEU:C	1:A:22:LEU:HD23	2.43	0.43
1:A:53:GLU:CD	7:A:612:HOH:O	2.61	0.43
1:A:44:SER:HA	1:A:142:SER:OG	2.18	0.43
1:A:258:HIS:CD2	1:B:108:MET:HE3	2.55	0.42
1:B:212:ALA:O	1:B:217:MET:CE	2.67	0.42
1:B:316:VAL:HA	1:B:329[A]:MET:HE3	1.98	0.42
1:B:46:VAL:O	1:B:85:VAL:HA	2.20	0.41
1:B:104:HIS:CE1	1:B:108:MET:HB2	2.56	0.41
1:B:140:LYS:NZ	7:B:628:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:LEU:C	1:B:274:LEU:HD23	2.46	0.41
1:A:73:TRP:CD1	1:A:184:ARG:HG2	2.56	0.41
1:B:198:THR:HA	1:B:236:MET:CE	2.50	0.41
1:A:244:ILE:HG12	1:A:256:TYR:O	2.20	0.41
1:A:388:ARG:CB	4:A:503:GOL:H31	2.48	0.41
1:B:435:LEU:HD23	1:B:440:LEU:HD22	2.03	0.41
1:A:94:THR:HG23	1:A:125:ASP:HB3	2.02	0.41
1:A:36:LYS:HB2	1:A:36:LYS:HE2	1.78	0.40
1:A:420:HIS:ND1	7:A:617:HOH:O	2.36	0.40
1:A:475:ILE:O	1:A:479:MET:HG3	2.21	0.40
1:A:140:LYS:N	1:A:141:PRO:CD	2.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:482:CYS:SG	1:B:482:CYS:SG[3_554]	1.53	0.67

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	490/488 (100%)	478 (98%)	12 (2%)	0	100	100
1	B	496/488 (102%)	488 (98%)	8 (2%)	0	100	100
All	All	986/976 (101%)	966 (98%)	20 (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	417/409 (102%)	417 (100%)	0	100	100
1	B	421/409 (103%)	420 (100%)	1 (0%)	87	84
All	All	838/818 (102%)	837 (100%)	1 (0%)	88	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	143	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	112	HIS
1	A	139	GLN
1	A	343	HIS
1	B	28	GLN
1	B	112	HIS
1	B	139	GLN

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

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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
5	GLY	B	503	2,3	3,3,4	0.80	0	1,2,4	0.25	0
5	GLY	A	504	2,3	3,3,4	0.76	0	1,2,4	0.28	0
4	GOL	A	503	-	5,5,5	0.18	0	5,5,5	0.31	0
6	PRO	A	505	-	8,8,8	0.84	1 (12%)	10,10,10	1.39	1 (10%)
6	PRO	B	504	-	8,8,8	0.83	1 (12%)	10,10,10	1.38	2 (20%)

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	GLY	B	503	2,3	-	0/0/1/2	-
5	GLY	A	504	2,3	-	0/0/1/2	-
4	GOL	A	503	-	-	0/4/4/4	-
6	PRO	A	505	-	-	0/4/11/11	0/1/1/1
6	PRO	B	504	-	-	0/4/11/11	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	505	PRO	OXT-C	-2.17	1.23	1.30
6	B	504	PRO	OXT-C	-2.13	1.23	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	505	PRO	OXT-C-O	-2.52	118.36	124.08
6	B	504	PRO	OXT-C-O	-2.47	118.48	124.08
6	B	504	PRO	OXT-C-CA	2.03	120.38	113.51

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	503	GOL	5	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	482/488 (98%)	-0.00	17 (3%)	47 51	14, 33, 68, 162	11 (2%)
1	B	479/488 (98%)	-0.02	17 (3%)	47 51	16, 33, 70, 125	18 (3%)
All	All	961/976 (98%)	-0.01	34 (3%)	47 51	14, 33, 69, 162	29 (3%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	489	PHE	10.4
1	B	484	LYS	7.0
1	A	488	PRO	6.7
1	A	486	PHE	5.9
1	B	241	TYR	5.3
1	A	487	THR	5.3
1	B	279	GLY	4.6
1	B	278	ASP	4.2
1	A	241	TYR	3.9
1	A	240	SER	3.8
1	B	239	SER	3.7
1	A	483	ASP	3.7
1	B	238	HIS	3.6
1	B	482	CYS	3.6
1	B	240	SER	3.5
1	B	168	LYS	3.3
1	A	278	ASP	3.3
1	A	343	HIS	3.0
1	A	482	CYS	2.9
1	B	483	ASP	2.8
1	B	243	CYS	2.7
1	A	279	GLY	2.6
1	B	260	GLY	2.6
1	A	275	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	92[A]	LYS	2.4
1	B	262	PRO	2.4
1	A	142	SER	2.4
1	B	275	PHE	2.3
1	B	480	ALA	2.3
1	A	481	GLY	2.3
1	A	168	LYS	2.3
1	A	259	ALA	2.2
1	B	159	ARG	2.2
1	B	437	ARG	2.1

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($\text{\AA}^2$)	Q<0.9
4	GOL	A	503	6/6	0.89	0.11	27,50,64,72	14
3	NA	B	502	1/1	0.94	0.07	36,36,36,36	0
5	GLY	A	504	4/5	0.94	0.09	23,37,44,44	0
5	GLY	B	503	4/5	0.95	0.09	26,44,55,55	0
3	NA	A	502	1/1	0.96	0.09	32,32,32,32	1
6	PRO	A	505	8/8	0.96	0.08	24,30,49,49	0
6	PRO	B	504	8/8	0.97	0.05	24,31,36,37	0
2	MN	B	501	1/1	0.99	0.06	25,25,25,25	1
2	MN	A	501	1/1	0.99	0.03	20,20,20,20	1

6.5 Other polymers [i](#)

There are no such residues in this entry.