



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 01:25 PM UTC

PDB ID : 2WUR / pdb_00002wur
Title : Atomic resolution structure of GFP measured on a rotating anode
Authors : Palm, G.J.; Schierbeek, A.J.; Kloos, M.
Deposited on : 2009-10-07
Resolution : 0.90 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

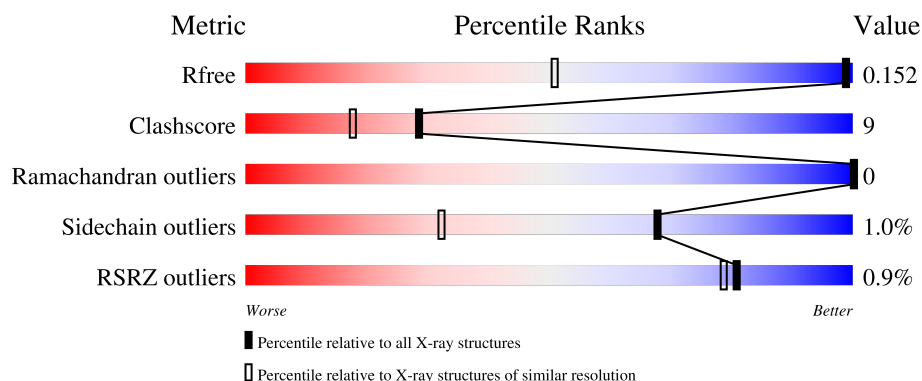
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 0.90 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	1366 (1.00-0.80)
Clashscore	190562	1416 (1.00-0.80)
Ramachandran outliers	187476	1341 (1.00-0.80)
Sidechain outliers	187428	1342 (1.00-0.80)
RSRZ outliers	180081	1363 (1.00-0.80)

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	236	

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
3	EOH	A	244	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EOH	A	245	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4296 atoms, of which 2004 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.

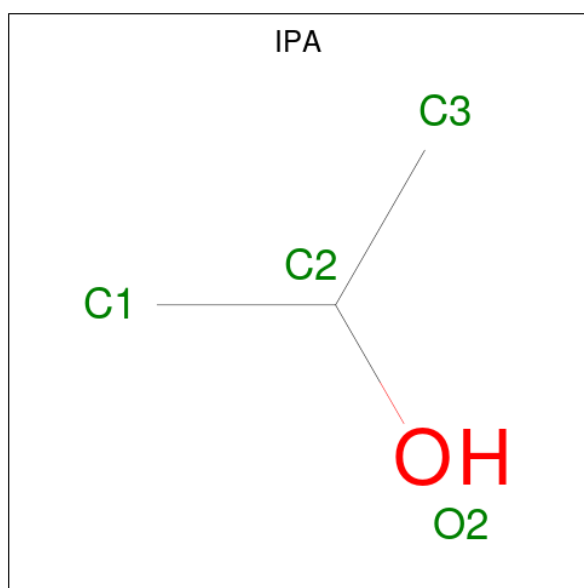
- Molecule 1 is a protein called GREEN FLUORESCENT PROTEIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	228	3885	1241	1943	328	366	7	0	22	1

There are 7 discrepancies between the modelled and reference sequences:

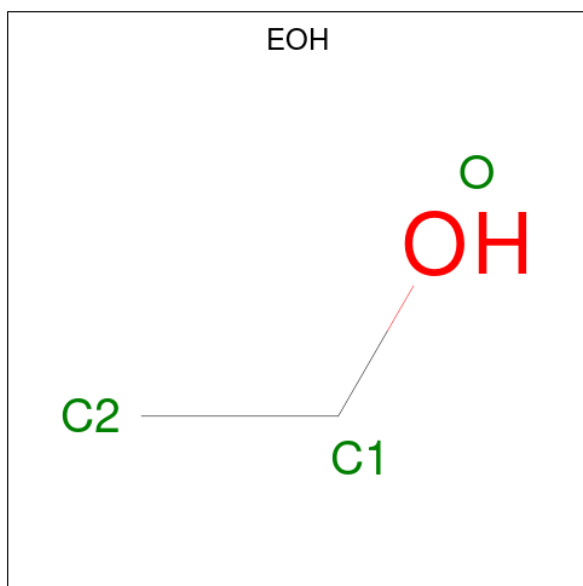
Chain	Residue	Modelled	Actual	Comment	Reference
A	64	LEU	PHE	engineered mutation	UNP P42212
A	66	GYS	SER	chromophore	UNP P42212
A	66	GYS	TYR	chromophore	UNP P42212
A	66	GYS	GLY	chromophore	UNP P42212
A	80	ARG	GLN	engineered mutation	UNP P42212
A	167	THR	ILE	engineered mutation	UNP P42212
A	238	ASN	LYS	engineered mutation	UNP P42212

- Molecule 2 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C_3H_8O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			12	3	8	1		

- Molecule 3 is ETHANOL (CCD ID: EOH) (formula: C_2H_6O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	1
			16	4	11	1		
3	A	1	Total	C	H	O	0	1
			18	4	12	2		
3	A	1	Total	C	H	O	0	0
			9	2	6	1		
3	A	1	Total	C	H	O	0	0
			9	2	6	1		
3	A	1	Total	C	H	O	0	0
			9	2	6	1		
3	A	1	Total	C	H	O	0	0
			9	2	6	1		

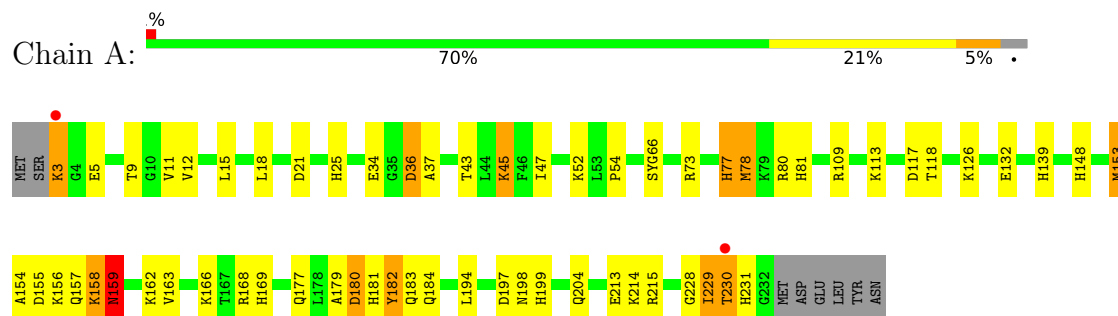
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	318	Total	O	0	2
			320	320		

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: GREEN FLUORESCENT PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.99Å 59.05Å 67.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 0.90 6.00 – 0.90	Depositor EDS
% Data completeness (in resolution range)	90.3 (6.00-0.90) 80.4 (6.00-0.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	-1.39 (at 0.90Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.146 , 0.174 0.129 , 0.152	Depositor DCC
R_{free} test set	13553 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	6.7	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.75 , 98.6	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	4296	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.88% 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: GYS, IPA, EOH

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	1.89	49/2026 (2.4%)	1.77	41/2732 (1.5%)

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	HIS	CE1-NE2	12.54	1.45	1.32
1	A	229	ILE	CA-C	-11.60	1.40	1.52
1	A	197	ASP	CG-OD2	-10.66	1.05	1.25
1	A	228	GLY	C-O	-10.48	1.12	1.23
1	A	168	ARG	NE-CZ	-9.18	1.23	1.33
1	A	231	HIS	C-N	-8.96	1.20	1.33
1	A	11	VAL	C-N	-8.80	1.26	1.33
1	A	132	GLU	CD-OE1	-8.17	1.09	1.25
1	A	213	GLU	CA-CB	8.13	1.65	1.53
1	A	231	HIS	CD2-NE2	8.08	1.46	1.37
1	A	36	ASP	CG-OD1	7.92	1.40	1.25
1	A	194	LEU	CA-CB	7.90	1.64	1.53
1	A	54	PRO	C-O	-7.80	1.14	1.24
1	A	3	LYS	N-CA	-7.61	1.33	1.49
1	A	156	LYS	C-N	-7.25	1.24	1.34
1	A	139	HIS	CG-ND1	-7.15	1.30	1.38
1	A	5	GLU	N-CA	-6.93	1.37	1.46
1	A	157	GLN	N-CA	6.91	1.54	1.46
1	A	25	HIS	CE1-NE2	6.86	1.39	1.32
1	A	231	HIS	CB-CG	-6.75	1.40	1.50
1	A	230	THR	C-O	6.54	1.32	1.23
1	A	228	GLY	C-N	-6.43	1.26	1.33
1	A	78[A]	MET	CA-CB	6.40	1.62	1.53
1	A	78[B]	MET	CA-CB	6.40	1.62	1.53
1	A	204	GLN	CG-CD	6.24	1.67	1.52
1	A	230	THR	CA-CB	6.21	1.63	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	230	THR	CA-C	-6.11	1.44	1.52
1	A	9	THR	N-CA	5.92	1.54	1.46
1	A	204	GLN	CD-OE1	-5.91	1.12	1.23
1	A	177	GLN	C-O	-5.90	1.16	1.24
1	A	11	VAL	CA-C	5.84	1.59	1.52
1	A	34	GLU	CD-OE2	-5.77	1.14	1.25
1	A	77	HIS	CD2-NE2	-5.77	1.31	1.37
1	A	179	ALA	C-O	-5.70	1.17	1.23
1	A	45	LYS	C-N	-5.67	1.25	1.33
1	A	229	ILE	C-O	5.57	1.30	1.24
1	A	194	LEU	CA-C	-5.52	1.46	1.52
1	A	81	HIS	CE1-NE2	5.47	1.38	1.32
1	A	154	ALA	CA-C	-5.45	1.45	1.52
1	A	37	ALA	C-N	-5.42	1.26	1.34
1	A	148	HIS	CE1-NE2	-5.35	1.27	1.32
1	A	113	LYS	CD-CE	-5.35	1.36	1.52
1	A	199	HIS	ND1-CE1	5.33	1.37	1.32
1	A	204	GLN	CB-CG	-5.27	1.36	1.52
1	A	153	MET	CA-CB	5.27	1.64	1.54
1	A	182	TYR	CE2-CZ	-5.14	1.25	1.38
1	A	169	HIS	C-O	-5.09	1.18	1.24
1	A	231	HIS	N-CA	-5.03	1.36	1.46
1	A	34	GLU	CB-CG	-5.01	1.37	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	ASP	CA-CB-CG	23.45	136.05	112.60
1	A	204	GLN	OE1-CD-NE2	21.39	143.99	122.60
1	A	229	ILE	CA-C-O	16.43	139.77	121.04
1	A	229	ILE	O-C-N	-15.85	105.40	122.67
1	A	168	ARG	CD-NE-CZ	15.45	146.03	124.40
1	A	73[A]	ARG	NE-CZ-NH1	-13.72	107.78	121.50
1	A	73[B]	ARG	NE-CZ-NH1	-13.72	107.78	121.50
1	A	168	ARG	NE-CZ-NH2	13.27	131.14	119.20
1	A	132	GLU	CB-CG-CD	10.73	130.84	112.60
1	A	197	ASP	CA-CB-CG	-10.65	101.95	112.60
1	A	80	ARG	NE-CZ-NH2	9.96	128.16	119.20
1	A	230	THR	OG1-CB-CG2	9.26	127.82	109.30
1	A	52	LYS	CB-CG-CD	9.10	132.24	111.30
1	A	168	ARG	NE-CZ-NH1	-9.07	112.43	121.50
1	A	204	GLN	CA-CB-CG	9.01	132.12	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	LYS	N-CA-CB	8.22	124.48	110.50
1	A	77	HIS	ND1-CE1-NE2	-8.19	100.21	108.40
1	A	157	GLN	OE1-CD-NE2	-8.10	114.50	122.60
1	A	153	MET	CG-SD-CE	-7.79	83.75	100.90
1	A	230	THR	N-CA-C	7.46	121.03	110.50
1	A	117	ASP	CA-CB-CG	7.13	119.73	112.60
1	A	204	GLN	CG-CD-NE2	-7.13	105.71	116.40
1	A	159	ASN	CA-CB-CG	-7.00	105.61	112.60
1	A	158	LYS	CG-CD-CE	-6.81	95.64	111.30
1	A	5	GLU	N-CA-CB	6.79	120.68	110.22
1	A	109	ARG	NE-CZ-NH2	-6.75	113.13	119.20
1	A	204	GLN	CG-CD-OE1	-6.52	107.75	120.80
1	A	109	ARG	NH1-CZ-NH2	6.30	127.50	119.30
1	A	157	GLN	CG-CD-OE1	6.25	133.29	120.80
1	A	156	LYS	CG-CD-CE	-6.24	96.94	111.30
1	A	180	ASP	CA-CB-CG	6.23	118.83	112.60
1	A	181	HIS	CE1-NE2-CD2	-6.13	102.87	109.00
1	A	159	ASN	CB-CG-OD1	-6.13	108.54	120.80
1	A	229	ILE	N-CA-C	-6.10	101.67	109.30
1	A	230	THR	O-C-N	-5.76	116.06	122.86
1	A	159	ASN	OD1-CG-ND2	5.72	128.32	122.60
1	A	12	VAL	CA-C-O	5.67	123.02	119.51
1	A	156	LYS	CA-C-O	-5.36	114.04	120.10
1	A	229	ILE	N-CA-CB	5.34	116.87	110.31
1	A	198	ASN	CA-CB-CG	-5.06	107.54	112.60
1	A	148	HIS	CE1-NE2-CD2	-5.01	103.99	109.00

There are no chirality outliers.

There are no planarity outliers.

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	1942	1943	1926	31	1
2	A	4	8	8	0	0
3	A	26	53	52	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	320	0	0	17	4
All	All	2292	2004	1986	35	5

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 9.

All (35) 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:229:ILE:HG23	4:A:2313:HOH:O	1.13	1.27
1:A:229:ILE:CG2	4:A:2313:HOH:O	1.65	1.25
1:A:3:LYS:HD2	4:A:2005:HOH:O	1.52	1.09
1:A:229:ILE:CA	4:A:2313:HOH:O	2.03	1.06
3:A:245:EOH:C1	4:A:2319:HOH:O	2.15	0.95
3:A:245:EOH:H11	4:A:2319:HOH:O	1.65	0.94
1:A:43[B]:THR:HG21	4:A:2076:HOH:O	1.74	0.87
1:A:230:THR:N	4:A:2313:HOH:O	2.07	0.86
3:A:245:EOH:C2	4:A:2319:HOH:O	2.35	0.75
1:A:36:ASP:OD2	4:A:2063:HOH:O	2.06	0.72
1:A:182:TYR:OH	4:A:2259:HOH:O	2.09	0.70
1:A:155:ASP:OD2	1:A:158:LYS:HD3	1.92	0.69
3:A:245:EOH:H21	4:A:2319:HOH:O	1.91	0.68
1:A:229:ILE:HA	4:A:2313:HOH:O	1.79	0.67
1:A:166[A]:LYS:HG2	1:A:180:ASP:OD1	1.97	0.64
1:A:15[B]:LEU:HG	1:A:118:THR:HG21	1.79	0.62
1:A:126:LYS:HD3	3:A:244:EOH:H21	1.85	0.59
1:A:126:LYS:CE	3:A:244:EOH:H21	2.35	0.56
1:A:158:LYS:HE2	1:A:184[B]:GLN:CD	2.31	0.55
1:A:3:LYS:O	4:A:2001:HOH:O	2.19	0.50
1:A:126:LYS:NZ	3:A:244:EOH:H21	2.28	0.48
1:A:126:LYS:CD	3:A:244:EOH:H21	2.42	0.48
1:A:3:LYS:HB2	4:A:2002:HOH:O	2.16	0.46
1:A:126:LYS:HD3	3:A:244:EOH:C2	2.45	0.46
1:A:214[B]:LYS:HE3	4:A:2320:HOH:O	2.15	0.45
1:A:126:LYS:NZ	3:A:244:EOH:C2	2.80	0.45
1:A:214[B]:LYS:HE2	4:A:2302:HOH:O	2.18	0.44
1:A:163:VAL:HB	1:A:183:GLN:HB3	2.00	0.44
1:A:77:HIS:CD2	1:A:78[B]:MET:HG2	2.54	0.43
1:A:45:LYS:HE2	1:A:47[B]:ILE:HD11	2.01	0.43
1:A:47[A]:ILE:HD13	1:A:215:ARG:CZ	2.51	0.41
1:A:21:ASP:OD1	1:A:21:ASP:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ASN:HD22	1:A:159:ASN:HA	1.71	0.41
1:A:78[B]:MET:HB3	1:A:78[B]:MET:HE2	1.69	0.40

All (5) 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:A:3:LYS:H2	1:A:162[B]:LYS:HZ2[3_645]	1.13	0.47
4:A:2198:HOH:O	4:A:2314:HOH:O[2_664]	1.77	0.43
4:A:2069:HOH:O	4:A:2271:HOH:O[3_655]	1.87	0.33
4:A:2116:HOH:O	4:A:2301:HOH:O[4_555]	2.06	0.14
4:A:2204:HOH:O	4:A:2276:HOH:O[2_664]	2.18	0.02

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	245/236 (104%)	242 (99%)	3 (1%)	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	220/206 (107%)	218 (99%)	2 (1%)	70	36

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	MET
1	A	159	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	135	ASN
1	A	170	ASN
1	A	177	GLN
1	A	212	ASN

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	GYS	A	66	1	20,22,23	1.90	5 (25%)	25,30,32	1.12	2 (8%)

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	GYS	A	66	1	-	1/10/29/30	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	GYS	CB2-CA2	4.66	1.39	1.35
1	A	66	GYS	CG2-CB2	-2.82	1.41	1.46
1	A	66	GYS	CA2-C2	-2.45	1.45	1.48
1	A	66	GYS	CA3-N3	-2.43	1.42	1.47
1	A	66	GYS	CE1-CD1	-2.43	1.34	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	GYS	CE2-CZ-CE1	-2.02	116.55	119.77
1	A	66	GYS	CD1-CE1-CZ	2.01	122.01	119.88

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	GYS	N1-CA1-CB1-OG1

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](#)

10 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
3	EOH	A	242[B]	-	2,2,2	0.41	0	1,1,1	0.22	0
3	EOH	A	241[A]	-	2,2,2	0.56	0	1,1,1	0.25	0
3	EOH	A	244	-	2,2,2	0.34	0	1,1,1	0.20	0
3	EOH	A	242[A]	-	2,2,2	0.36	0	1,1,1	0.09	0
3	EOH	A	245	-	2,2,2	0.54	0	1,1,1	0.30	0
3	EOH	A	247	-	2,2,2	0.43	0	1,1,1	0.32	0
3	EOH	A	246	-	2,2,2	0.51	0	1,1,1	0.30	0
3	EOH	A	243	-	2,2,2	0.40	0	1,1,1	0.21	0
2	IPA	A	240	-	3,3,3	2.15	1 (33%)	3,3,3	1.26	1 (33%)
3	EOH	A	241[B]	-	2,2,2	0.49	0	1,1,1	0.30	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	240	IPA	C1-C2	3.42	1.71	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	240	IPA	C3-C2-C1	-2.14	97.36	113.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	244	EOH	6	0
3	A	245	EOH	4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	227/236 (96%)	-0.93	2 (0%) 81 79	4, 8, 17, 43	25 (11%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	230	THR	2.7
1	A	3	LYS	2.3

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	GYS	A	66	21/22	1.00	0.02	3,4,5,8	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($\text{\AA}^2$)	Q<0.9
3	EOH	A	243	3/3	0.79	0.10	32,39,64,64	0
3	EOH	A	247	3/3	0.80	0.10	18,27,41,43	0
2	IPA	A	240	4/4	0.81	0.15	8,24,30,31	0
3	EOH	A	241[A]	3/3	0.89	0.13	1,14,30,30	7
3	EOH	A	241[B]	3/3	0.89	0.13	3,8,14,21	7
3	EOH	A	244	3/3	0.92	0.10	4,9,18,27	0
3	EOH	A	245	3/3	0.95	0.09	9,14,34,34	0
3	EOH	A	242[A]	3/3	0.96	0.09	9,12,27,41	9
3	EOH	A	242[B]	3/3	0.96	0.09	10,16,24,29	9
3	EOH	A	246	3/3	0.98	0.06	4,14,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.