



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:09 PM UTC

PDB ID : 8OVN / pdb_00008ovn
Title : X-ray structure of the SF-iGluSnFR-S72A
Authors : Tarnawski, M.; Hellweg, L.; Bergner, A.; Hiblot, J.; Leippe, P.; Johnsson, K.
Deposited on : 2023-04-26
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

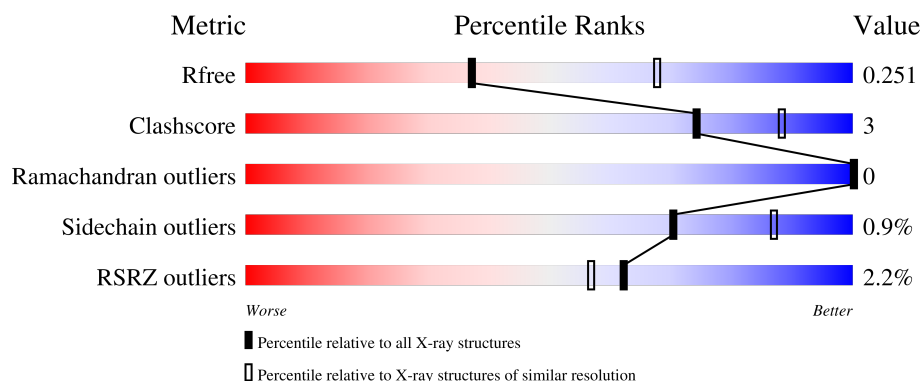
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	518	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4019 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 Putative periplasmic binding transport protein, Green fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	504	Total	C	N	O	S	0	0	0
			3987	2518	682	771	16			

There are 31 discrepancies between the modelled and reference sequences:

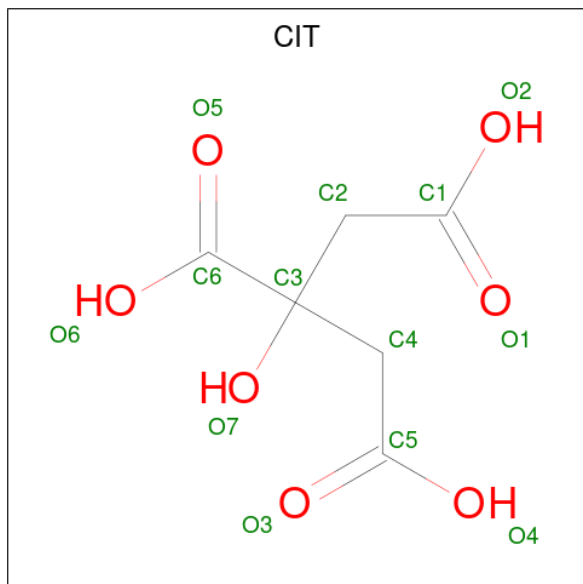
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP A0A0H2UXX1
A	69	ALA	SER	engineered mutation	UNP A0A0H2UXX1
A	124	ASN	ASP	engineered mutation	UNP A0A0H2UXX1
A	127	ASP	GLY	engineered mutation	UNP A0A0H2UXX1
A	181	VAL	ALA	engineered mutation	UNP A0A0H2UXX1
A	197	GLU	ASP	engineered mutation	UNP A0A0H2UXX1
A	249	LEU	PRO	engineered mutation	UNP A0A0H2UXX1
A	250	VAL	PRO	engineered mutation	UNP A0A0H2UXX1
A	257	THR	MET	engineered mutation	UNP P42212
A	267	ALA	VAL	engineered mutation	UNP P42212
A	275	VAL	ILE	engineered mutation	UNP P42212
A	310	VAL	ALA	engineered mutation	UNP P42212
A	335	LEU	HIS	engineered mutation	UNP P42212
A	343	GLY	-	linker	UNP P42212
A	344	GLY	-	linker	UNP P42212
A	345	THR	-	linker	UNP P42212
A	346	GLY	-	linker	UNP P42212
A	347	GLY	-	linker	UNP P42212
A	348	SER	-	linker	UNP P42212
A	378	ARG	SER	engineered mutation	UNP P42212
A	387	ASN	TYR	engineered mutation	UNP P42212
A	412	LEU	PHE	engineered mutation	UNP P42212
A	413	CRO	SER	chromophore	UNP P42212
A	413	CRO	TYR	chromophore	UNP P42212
A	413	CRO	GLY	chromophore	UNP P42212
A	447	SER	PHE	engineered mutation	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	453	THR	ASN	engineered mutation	UNP P42212
A	493	PHE	TYR	engineered mutation	UNP P42212
A	495	ASN	SER	engineered mutation	UNP P42212
A	496	PRO	HIS	engineered mutation	UNP P42212
A	520	LYS	ASN	engineered mutation	UNP A0A0H2UXX1

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 13 6 7	0	0

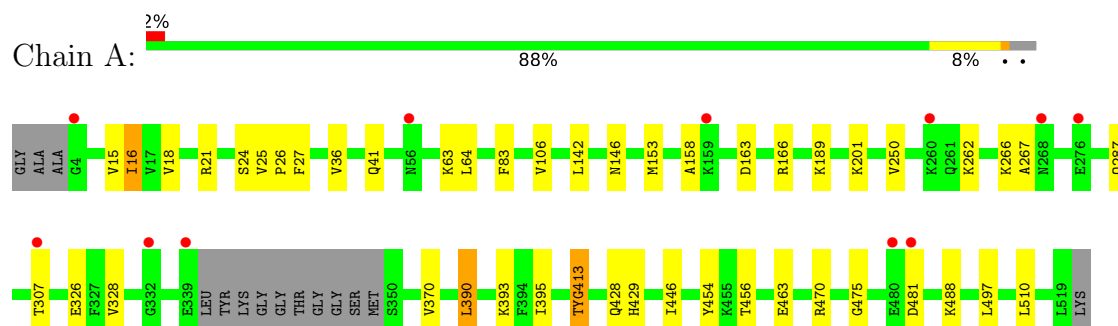
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	19	Total O 19 19	0	0

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: Putative periplasmic binding transport protein, Green fluorescent protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.14Å 106.14Å 108.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.60 – 2.60 46.60 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.60-2.60) 99.9 (46.60-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.200 , 0.253 0.199 , 0.251	Depositor DCC
R_{free} test set	1105 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4019	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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: CIT, CRO

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	0.08	0/4038	0.24	0/5445

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

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	3987	0	3953	25	0
2	A	13	0	5	1	0
3	A	19	0	0	0	0
All	All	4019	0	3958	25	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 3.

All (25) 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:326:GLU:HB3	1:A:390:LEU:HB2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:THR:HG23	1:A:328:VAL:HG22	1.77	0.66
1:A:267:ALA:HB3	1:A:287:GLN:HB3	1.78	0.65
1:A:146:ASN:HB2	1:A:153:MET:HE2	1.82	0.62
1:A:446:ILE:HB	1:A:454:TYR:HB2	1.86	0.58
1:A:21:ARG:NH1	2:A:601:CIT:O2	2.37	0.55
1:A:15:VAL:HG11	1:A:63:LYS:HD2	1.90	0.53
1:A:428:GLN:HG2	1:A:429:HIS:CD2	2.48	0.49
1:A:370:VAL:HG22	1:A:475:GLY:HA3	1.96	0.47
1:A:158:ALA:HB1	1:A:163:ASP:HB3	1.96	0.47
1:A:189:LYS:HG2	1:A:497:LEU:HD21	1.95	0.47
1:A:142:LEU:HG	1:A:153:MET:HE1	1.98	0.46
1:A:163:ASP:OD1	1:A:166:ARG:NH2	2.46	0.46
1:A:262:LYS:NZ	1:A:266:LYS:HZ1	2.15	0.45
1:A:18:VAL:HB	1:A:64:LEU:HD23	2.00	0.45
1:A:36:VAL:HG12	1:A:41:GLN:HG2	1.99	0.45
1:A:481:ASP:O	1:A:488:LYS:NZ	2.46	0.44
1:A:463:GLU:OE2	1:A:470:ARG:NH1	2.51	0.44
1:A:25:VAL:HA	1:A:26:PRO:HA	1.80	0.43
1:A:24:SER:O	1:A:27:PHE:N	2.50	0.43
1:A:106:VAL:HG23	1:A:510:LEU:HD22	2.01	0.43
1:A:16:ILE:HD11	1:A:83:PHE:HE1	1.85	0.42
1:A:413:CRO:N2	1:A:413:CRO:HD1	2.36	0.41
1:A:201:LYS:HA	1:A:201:LYS:HD3	1.77	0.41
1:A:393:LYS:HE2	1:A:395:ILE:HD11	2.02	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	497/518 (96%)	486 (98%)	11 (2%)	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	439/446 (98%)	435 (99%)	4 (1%)	70 87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	250	VAL
1	A	390	LEU
1	A	456	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	152	ASN
1	A	229	GLN
1	A	253	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	CRO	A	413	1	22,23,24	0.41	0	30,32,34	0.73	1 (3%)

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	CRO	A	413	1	-	1/12/31/32	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	CRO	N3-C1-N2	2.58	113.51	111.48

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	413	CRO	C3-CA3-N3-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	413	CRO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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	$\# Z > 2$
2	CIT	A	601	-	12,12,12	1.02	0	17,17,17	2.16	8 (47%)

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	CIT	A	601	-	-	8/16/16/16	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	CIT	C4-C3-C6	5.25	121.64	110.03
2	A	601	CIT	C3-C4-C5	3.06	122.28	113.92
2	A	601	CIT	O6-C6-C3	2.88	118.66	113.14
2	A	601	CIT	O7-C3-C6	-2.45	105.49	108.96
2	A	601	CIT	O2-C1-O1	-2.27	117.49	123.33
2	A	601	CIT	O6-C6-O5	-2.14	117.02	123.86
2	A	601	CIT	O4-C5-O3	-2.07	118.02	123.33
2	A	601	CIT	O7-C3-C4	-2.06	104.68	109.38

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	CIT	O7-C3-C4-C5
2	A	601	CIT	O7-C3-C6-O5
2	A	601	CIT	O7-C3-C6-O6
2	A	601	CIT	C4-C3-C6-O5
2	A	601	CIT	C4-C3-C6-O6
2	A	601	CIT	C2-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
2	A	601	CIT	C6-C3-C4-C5
2	A	601	CIT	C1-C2-C3-O7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	CIT	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 [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	503/518 (97%)	-0.05	11 (2%) 62 57	34, 46, 68, 89	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	56	ASN	3.0
1	A	332	GLY	2.9
1	A	480	GLU	2.4
1	A	159	LYS	2.4
1	A	481	ASP	2.3
1	A	276	GLU	2.2
1	A	339	GLU	2.1
1	A	260	LYS	2.1
1	A	307	THR	2.0
1	A	268	ASN	2.0
1	A	4	GLY	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	CRO	A	413	22/23	0.94	0.09	40,47,55,61	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

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
2	CIT	A	601	13/13	0.79	0.14	44,57,68,69	0

6.5 Other polymers

There are no such residues in this entry.