wwPDB X-ray Structure Validation Summary Report (i)

Mar 14, 2026 – 10:09 AM UTC

PDB ID	:	5IZS / pdb_00005izs
Title	:	De novo design of protein homo-oligomers with modular hydrogen bond network-mediated specificity
Authors	:	Sankaran, B.; Zwart, P.H.; Pereira, J.H.; Baker, D.; Boyken, S.; Chen, Z.; Oberdorfer, G.
Deposited on	:	2016-03-25
Resolution	:	2.36 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

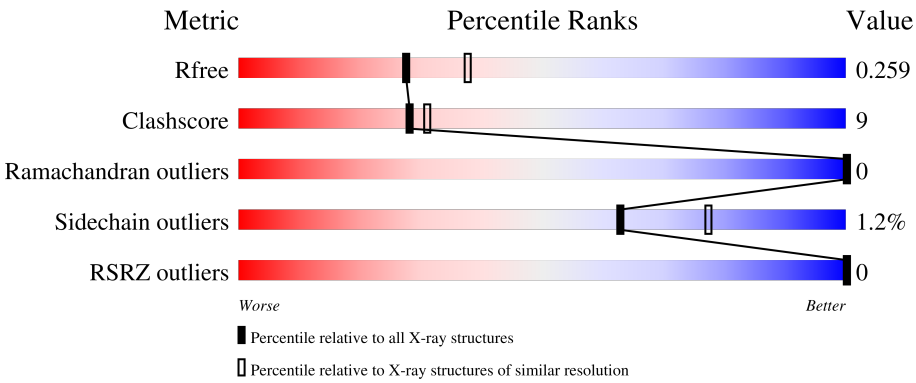
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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 <=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	95	67% 14% 19%
1	B	95	61% 20% 19%
1	C	95	77% 9% 14%
1	D	95	69% 16% • 14%
1	E	95	59% 25% • 15%

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Mol	Chain	Length	Quality of chain
1	F	95	71% 14% 16%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4127 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 Designed protein 5L6HC3_1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	77	637	391	118	128	0	0	0
1	B	77	648	397	122	129	0	1	0
1	C	82	681	416	129	136	0	0	0
1	D	82	680	416	126	138	0	0	0
1	E	81	691	422	130	139	0	2	0
1	F	80	665	408	124	133	0	0	0

- Molecule 2 is water.

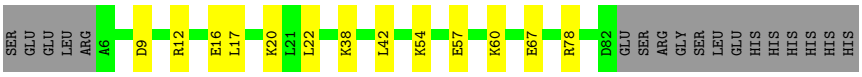
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	24	24	24	0	0
2	B	23	23	23	0	0
2	C	22	22	22	0	0
2	D	17	17	17	0	0
2	E	22	22	22	0	0
2	F	17	17	17	0	0

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.

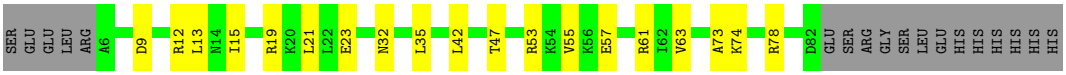
• Molecule 1: Designed protein 5L6HC3_1

Chain A: 67% 14% 19%



• Molecule 1: Designed protein 5L6HC3_1

Chain B: 61% 20% 19%

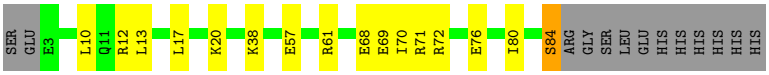


• Molecule 1: Designed protein 5L6HC3_1

Chain C: 77% 9% 14%

• Molecule 1: Designed protein 5L6HC3_1

Chain D: 69% 16% • 14%

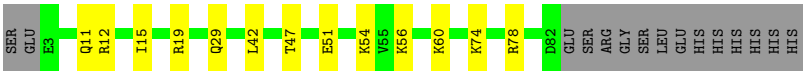


• Molecule 1: Designed protein 5L6HC3_1

Chain E: 59% 25% • 15%

• Molecule 1: Designed protein 5L6HC3_1

Chain F: 71% 14% 16%


SER
GLU
E3
Q11
R12
I15
R19
Q29
L42
T47
E51
K54
V55
K56
K60
K74
R78
D82
GLU
SER
ARG
GLY
SER
LEU
GLU
HIS
HIS
HIS
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4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α, β, γ	72.67Å 72.67Å 88.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	51.24 – 2.36 51.24 – 2.36	Depositor EDS
% Data completeness (in resolution range)	99.0 (51.24-2.36) 99.1 (51.24-2.36)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.44 (at 2.37Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.245 , 0.269 0.247 , 0.259	Depositor DCC
R_{free} test set	2033 reflections (9.56%)	wwPDB-VP
Wilson B-factor (Å^2)	22.4	Xtriage
Anisotropy	0.187	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 35.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.055 for -h,-k,l 0.068 for h,-h-k,-l 0.220 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4127	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.03 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.3676e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.30	0/637	0.77	0/850
1	B	0.29	0/648	0.71	0/864
1	C	0.32	0/681	0.74	0/907
1	D	0.29	0/680	0.76	0/907
1	E	0.33	0/691	0.76	0/921
1	F	0.29	0/665	0.73	0/887
All	All	0.31	0/4002	0.74	0/5336

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	637	0	666	11	0
1	B	648	0	678	19	0
1	C	681	0	710	10	0
1	D	680	0	707	16	0
1	E	691	0	718	24	0
1	F	665	0	696	9	0
2	A	24	0	0	1	0
2	B	23	0	0	5	0
2	C	22	0	0	0	0
2	D	17	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	22	0	0	6	0
2	F	17	0	0	0	0
All	All	4127	0	4175	73	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 9.

The worst 5 of 73 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:C:57:GLU:OE2	1:C:61:ARG:NH1	1.95	0.99
1:E:17:LEU:HA	1:E:20:LYS:HE3	1.57	0.86
1:E:11:GLN:NE2	2:E:102:HOH:O	2.12	0.82
1:F:51:GLU:OE1	1:F:54:LYS:NZ	2.12	0.82
1:E:71:ARG:NH1	2:E:101:HOH:O	2.05	0.81

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	75/95 (79%)	75 (100%)	0	0	100	100
1	B	76/95 (80%)	76 (100%)	0	0	100	100
1	C	80/95 (84%)	80 (100%)	0	0	100	100
1	D	80/95 (84%)	79 (99%)	1 (1%)	0	100	100
1	E	81/95 (85%)	80 (99%)	1 (1%)	0	100	100
1	F	78/95 (82%)	78 (100%)	0	0	100	100
All	All	470/570 (82%)	468 (100%)	2 (0%)	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	69/86 (80%)	68 (99%)	1 (1%)	59	73
1	B	70/86 (81%)	70 (100%)	0	100	100
1	C	74/86 (86%)	73 (99%)	1 (1%)	59	73
1	D	74/86 (86%)	73 (99%)	1 (1%)	59	73
1	E	75/86 (87%)	73 (97%)	2 (3%)	39	52
1	F	72/86 (84%)	72 (100%)	0	100	100
All	All	434/516 (84%)	429 (99%)	5 (1%)	63	77

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	GLU
1	C	84	SER
1	D	84	SER
1	E	55	VAL
1	E	65	GLU

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

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 (i)

There are no ligands in this entry.

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	77/95 (81%)	-1.22	0	100	100	20, 37, 54, 66	0
1	B	77/95 (81%)	-1.23	0	100	100	18, 34, 54, 66	1 (1%)
1	C	82/95 (86%)	-1.20	0	100	100	22, 37, 61, 80	0
1	D	82/95 (86%)	-1.28	0	100	100	19, 37, 54, 63	0
1	E	81/95 (85%)	-1.24	0	100	100	16, 38, 55, 75	2 (2%)
1	F	80/95 (84%)	-1.25	0	100	100	22, 35, 52, 64	0
All	All	479/570 (84%)	-1.23	0	100	100	16, 37, 57, 80	3 (0%)

There are no RSRZ outliers to report.

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.