wwPDB X-ray Structure Validation Summary Report (i)

Sep 12, 2023 – 09:41 AM EDT

PDB ID	:	4NN7
Title	:	Cytokine receptor complex - Crystal form 2
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Deposited on	:	2013-11-16
Resolution	:	3.77 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 (i)) were used in the production of this report:

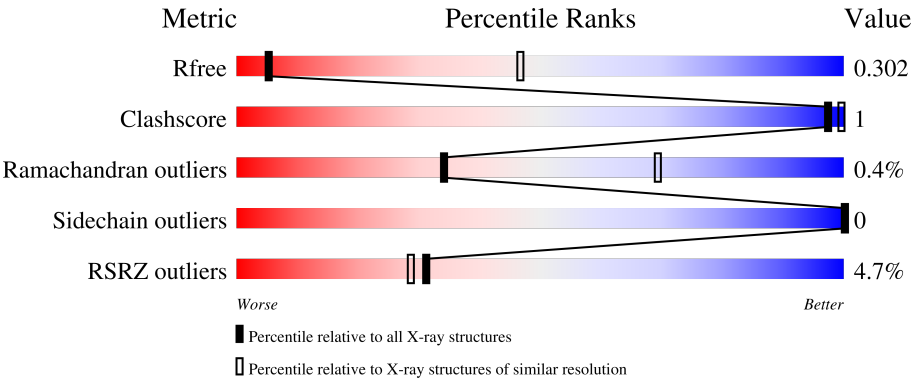
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 3.77 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}	130704	1038 (3.96-3.60)
Clashscore	141614	1100 (3.96-3.60)
Ramachandran outliers	138981	1062 (3.96-3.60)
Sidechain outliers	138945	1058 (3.96-3.60)
RSRZ outliers	127900	1009 (3.98-3.58)

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	130	2% 78% 5% 17%
2	B	223	6% 87% 13%
3	C	212	3% 78% • 20%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6114 atoms, of which 2817 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 Thymic stromal lymphopoietin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	108	Total	C	H	N	O	S	0	0	0
			1435	498	662	125	142	8			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	GLN	ASN	engineered mutation	UNP Q9JIE6
A	141	GLY	-	expression tag	UNP Q9JIE6
A	142	THR	-	expression tag	UNP Q9JIE6
A	143	LYS	-	expression tag	UNP Q9JIE6
A	144	HIS	-	expression tag	UNP Q9JIE6
A	145	HIS	-	expression tag	UNP Q9JIE6
A	146	HIS	-	expression tag	UNP Q9JIE6
A	147	HIS	-	expression tag	UNP Q9JIE6
A	148	HIS	-	expression tag	UNP Q9JIE6
A	149	HIS	-	expression tag	UNP Q9JIE6

- Molecule 2 is a protein called Interleukin-7 receptor subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	195	Total	C	H	N	O	S	0	0	0
			2421	858	1096	224	236	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	17	GLY	-	expression tag	UNP P16872
B	18	SER	-	expression tag	UNP P16872
B	19	HIS	-	expression tag	UNP P16872
B	20	MET	-	expression tag	UNP P16872

- Molecule 3 is a protein called Cytokine receptor-like factor 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	170	Total	C	H	N	O	S	0	0	0
			2258	765	1059	205	222	7			

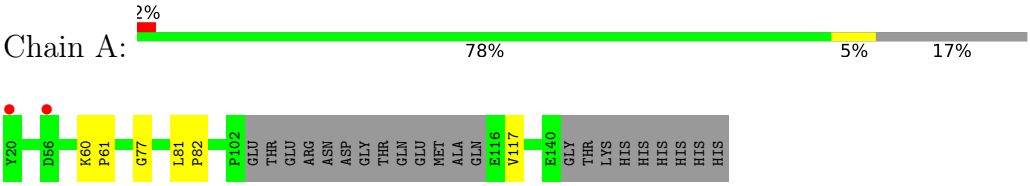
There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	53	GLN	ASN	engineered mutation	UNP Q8CII9
C	179	VAL	ALA	conflict	UNP Q8CII9
C	223	GLY	-	expression tag	UNP Q8CII9
C	224	THR	-	expression tag	UNP Q8CII9
C	225	LYS	-	expression tag	UNP Q8CII9
C	226	HIS	-	expression tag	UNP Q8CII9
C	227	HIS	-	expression tag	UNP Q8CII9
C	228	HIS	-	expression tag	UNP Q8CII9
C	229	HIS	-	expression tag	UNP Q8CII9
C	230	HIS	-	expression tag	UNP Q8CII9
C	231	HIS	-	expression tag	UNP Q8CII9

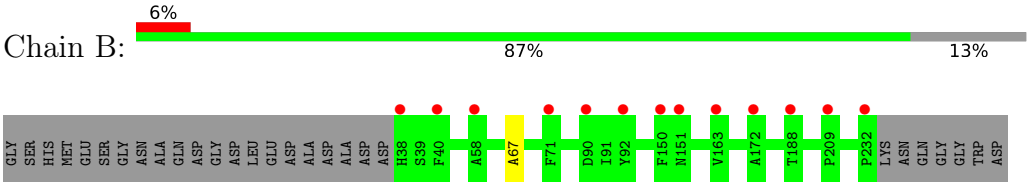
3 Residue-property plots (i)

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.

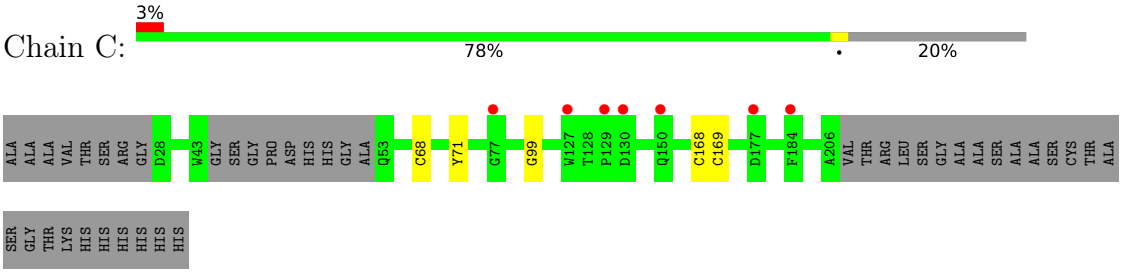
• Molecule 1: Thymic stromal lymphopoietin



• Molecule 2: Interleukin-7 receptor subunit alpha



• Molecule 3: Cytokine receptor-like factor 2



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	35.79Å 50.10Å 249.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.13 – 3.77 49.13 – 3.77	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.13-3.77) 99.5 (49.13-3.77)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.95 (at 3.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.3_1479)	Depositor
R, R_{free}	0.277 , 0.286 0.289 , 0.302	Depositor DCC
R_{free} test set	249 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	98.8	Xtriage
Anisotropy	0.865	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 93.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.39$, $< L^2 > = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.86	EDS
Total number of atoms	6114	wwPDB-VP
Average B, all atoms (Å^2)	109.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.30% of the height of the origin peak. No significant pseudotranslation is detected.*

[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

5.1 Standard geometry

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	0.27	0/794	0.42	0/1090
2	B	0.25	0/1361	0.44	0/1886
3	C	0.25	0/1235	0.45	0/1703
All	All	0.26	0/3390	0.44	0/4679

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	773	662	660	3	0
2	B	1325	1096	1091	0	0
3	C	1199	1059	1055	2	0
All	All	3297	2817	2806	5	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 1.

All (5) 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)
3:C:68:CYS:HB3	3:C:71:TYR:CE2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:GLY:O	1:A:117:VAL:N	2.45	0.47
1:A:81:LEU:HB3	1:A:82:PRO:HD3	1.98	0.44
1:A:60:LYS:N	1:A:61:PRO:HD2	2.32	0.44
3:C:168:CYS:HA	3:C:169:CYS:HA	1.87	0.42

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	104/130 (80%)	103 (99%)	1 (1%)	0	100	100
2	B	193/223 (86%)	184 (95%)	8 (4%)	1 (0%)	29	65
3	C	166/212 (78%)	164 (99%)	1 (1%)	1 (1%)	25	61
All	All	463/565 (82%)	451 (97%)	10 (2%)	2 (0%)	34	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	67	ALA
3	C	99	GLY

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	73/119 (61%)	73 (100%)	0	100	100
2	B	108/198 (54%)	108 (100%)	0	100	100
3	C	108/165 (66%)	108 (100%)	0	100	100
All	All	289/482 (60%)	289 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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 monosaccharides 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

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, 95^{th} 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	108/130 (83%)	0.12	2 (1%)	66	63	71, 89, 118, 137	0
2	B	195/223 (87%)	0.46	13 (6%)	17	15	76, 109, 152, 220	0
3	C	170/212 (80%)	0.43	7 (4%)	37	33	68, 90, 125, 146	0
All	All	473/565 (83%)	0.37	22 (4%)	31	29	68, 98, 141, 220	0

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	163	VAL	3.6
2	B	71	PHE	3.4
2	B	90	ASP	3.4
2	B	172	ALA	3.4
3	C	130	ASP	3.4

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.