



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 12:14 PM EDT

PDB ID : 7AX1
Title : Crystal structure of the human CCR4-CAF1 complex
Authors : Chen, Y.; Khazina, E.; Weichenrieder, O.
Deposited on : 2020-11-09
Resolution : 3.30 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	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

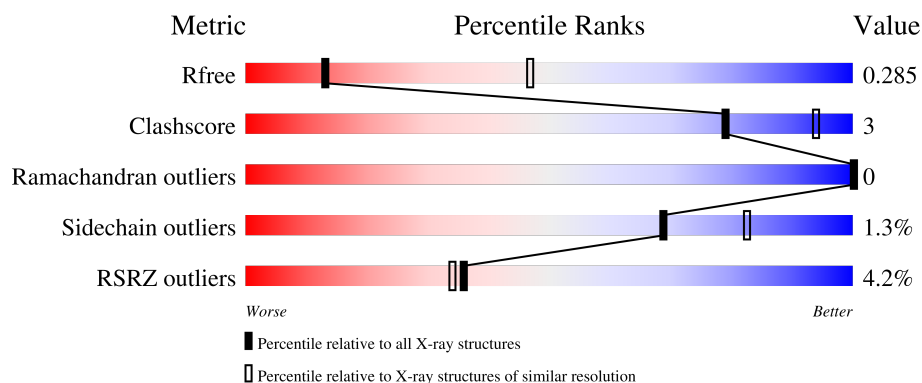
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	563	<div> <div>4%</div> <div>85%</div> <div>6% • 8%</div> </div>
2	B	290	<div> <div>3%</div> <div>80%</div> <div>8% 13%</div> </div>

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	MG	A	1001	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12393 atoms, of which 6160 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 CCR4-NOT transcription complex subunit 6.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	518	8301	2653	4143	706	778	21	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP Q9ULM6
A	-4	PRO	-	expression tag	UNP Q9ULM6
A	-3	HIS	-	expression tag	UNP Q9ULM6
A	-2	MET	-	expression tag	UNP Q9ULM6
A	-1	LEU	-	expression tag	UNP Q9ULM6
A	0	GLU	-	expression tag	UNP Q9ULM6

- Molecule 2 is a protein called CCR4-NOT transcription complex subunit 7.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	253	4090	1338	2017	332	388	15	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP Q9UIV1
B	-3	PRO	-	expression tag	UNP Q9UIV1
B	-2	HIS	-	expression tag	UNP Q9UIV1
B	-1	MET	-	expression tag	UNP Q9UIV1
B	0	LEU	-	expression tag	UNP Q9UIV1
B	1	GLU	-	expression tag	UNP Q9UIV1

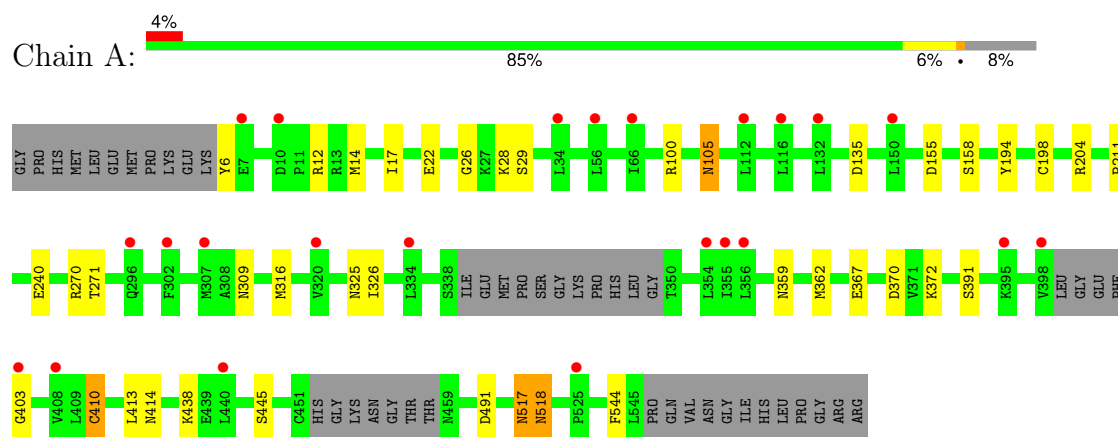
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

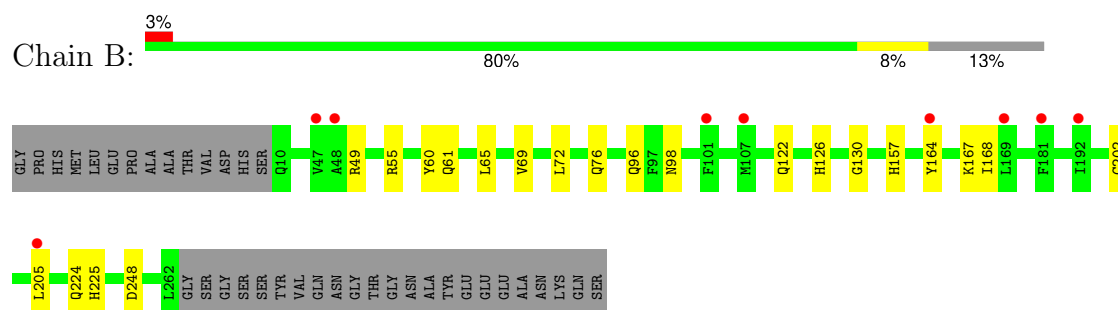
3 Residue-property plots [i](#)

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: CCR4-NOT transcription complex subunit 6



- Molecule 2: CCR4-NOT transcription complex subunit 7



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	149.85Å 63.26Å 113.07Å 90.00° 106.81° 90.00°	Depositor
Resolution (Å)	71.72 – 3.30 71.72 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (71.72-3.30) 99.8 (71.72-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.248 , 0.276 0.259 , 0.285	Depositor DCC
R_{free} test set	776 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	150.5	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 104.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12393	wwPDB-VP
Average B, all atoms (Å ²)	163.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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: MG

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	0/4250	0.43	0/5757
2	B	0.26	0/2121	0.39	0/2859
All	All	0.27	0/6371	0.42	0/8616

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	4158	4143	4143	24	0
2	B	2073	2017	2017	15	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	6233	6160	6160	35	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 (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)
2:B:224:GLN:OE1	2:B:225:HIS:ND1	2.33	0.62
1:A:22:GLU:O	1:A:26:GLY:N	2.28	0.62
1:A:367:GLU:O	1:A:438:LYS:NZ	2.32	0.61
2:B:72:LEU:O	2:B:167:LYS:NZ	2.36	0.58
2:B:69:VAL:O	2:B:167:LYS:NZ	2.28	0.57
1:A:12:ARG:O	2:B:49:ARG:NH1	2.36	0.56
2:B:76:GLN:NE2	2:B:96:GLN:OE1	2.40	0.55
1:A:316:MET:SD	1:A:445:SER:OG	2.58	0.55
1:A:325:ASN:OD1	1:A:326:ILE:N	2.41	0.54
1:A:204:ARG:NH1	1:A:211:PRO:HA	2.24	0.53
1:A:22:GLU:OE2	2:B:55:ARG:HA	2.08	0.53
1:A:391:SER:OG	1:A:403:GLY:O	2.26	0.53
2:B:122:GLN:O	2:B:126:HIS:ND1	2.38	0.53
1:A:29:SER:OG	2:B:55:ARG:N	2.44	0.50
2:B:164:TYR:O	2:B:168:ILE:HG12	2.13	0.49
1:A:362:MET:HG2	1:A:413:LEU:HA	1.93	0.49
1:A:370:ASP:N	1:A:370:ASP:OD1	2.48	0.47
1:A:17:ILE:HD12	1:A:22:GLU:OE2	2.15	0.47
1:A:517:ASN:N	1:A:517:ASN:OD1	2.48	0.46
1:A:270:ARG:O	1:A:271:THR:OG1	2.30	0.46
1:A:309:ASN:OD1	1:A:309:ASN:N	2.49	0.45
1:A:14:MET:HE1	2:B:60:TYR:CE2	2.52	0.44
1:A:105:ASN:N	1:A:105:ASN:OD1	2.50	0.44
1:A:6:TYR:OH	1:A:100:ARG:HB3	2.18	0.43
1:A:359:ASN:HA	1:A:410:CYS:O	2.19	0.43
2:B:76:GLN:NE2	2:B:98:ASN:OD1	2.45	0.43
2:B:61:GLN:O	2:B:65:LEU:HD13	2.19	0.43
2:B:126:HIS:O	2:B:130:GLY:N	2.51	0.43
1:A:155:ASP:O	1:A:158:SER:OG	2.32	0.42
1:A:28:LYS:HD2	1:A:28:LYS:H	1.84	0.42
1:A:198:CYS:H	1:A:240:GLU:HB3	1.85	0.42
2:B:202:CYS:HB2	2:B:205:LEU:HD12	2.01	0.41
2:B:157:HIS:O	2:B:157:HIS:ND1	2.53	0.41
1:A:518:ASN:OD1	1:A:518:ASN:N	2.54	0.41
1:A:372:LYS:NZ	1:A:414:ASN:O	2.42	0.40

There are no symmetry-related clashes.

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	510/563 (91%)	477 (94%)	33 (6%)	0	100	100
2	B	251/290 (87%)	241 (96%)	10 (4%)	0	100	100
All	All	761/853 (89%)	718 (94%)	43 (6%)	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	468/505 (93%)	460 (98%)	8 (2%)	60	78
2	B	225/253 (89%)	224 (100%)	1 (0%)	91	95
All	All	693/758 (91%)	684 (99%)	9 (1%)	69	82

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	135	ASP
1	A	194	TYR
1	A	410	CYS
1	A	491	ASP
1	A	517	ASN
1	A	518	ASN
1	A	544	PHE

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Mol	Chain	Res	Type
2	B	248	ASP

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	518/563 (92%)	0.35	23 (4%) 34 33	82, 135, 220, 279	0
2	B	253/290 (87%)	0.33	9 (3%) 42 40	122, 157, 207, 244	0
All	All	771/853 (90%)	0.35	32 (4%) 36 34	82, 145, 217, 279	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	101	PHE	4.0
1	A	56	LEU	3.9
1	A	302	PHE	3.3
1	A	7	GLU	3.1
2	B	164	TYR	2.9
1	A	403	GLY	2.9
1	A	356	LEU	2.9
1	A	408	VAL	2.9
2	B	192	ILE	2.8
2	B	181	PHE	2.8
1	A	112	LEU	2.7
2	B	169	LEU	2.7
2	B	205	LEU	2.6
1	A	10	ASP	2.5
1	A	132	LEU	2.5
1	A	34	LEU	2.5
1	A	395	LYS	2.5
1	A	440	LEU	2.4
1	A	398	VAL	2.4
2	B	47	VAL	2.3
2	B	107	MET	2.3
1	A	116	LEU	2.3
2	B	48	ALA	2.2
1	A	320	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	307	MET	2.2
1	A	525	PRO	2.1
1	A	150	LEU	2.1
1	A	334	LEU	2.1
1	A	355	ILE	2.0
1	A	296	GLN	2.0
1	A	66	ILE	2.0
1	A	354	LEU	2.0

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 monosaccharides 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(\AA^2)	Q<0.9
3	MG	B	1001	1/1	0.48	0.32	158,158,158,158	0
3	MG	A	1001	1/1	0.77	0.68	121,121,121,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.