



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8D30
Title : Crystal structure of the human COPB2 WD-domains
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Deposited on : 2022-05-31
Resolution : 2.40 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
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
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)	:	2.31.2

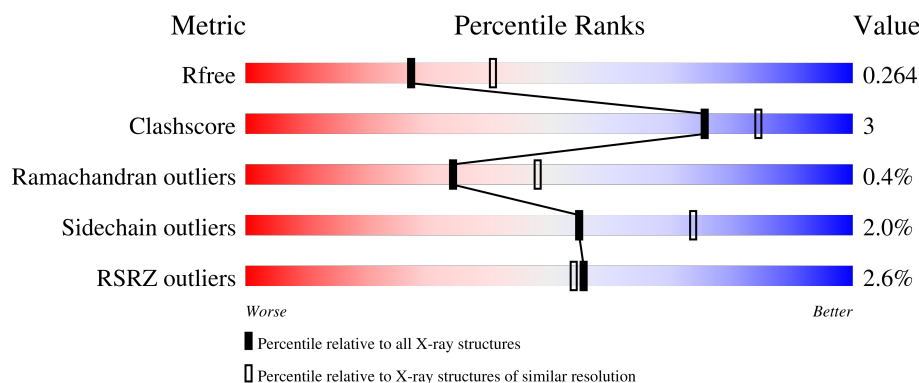
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	615	
1	B	615	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8563 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 Coatomer subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	586	Total	C	N	O	S	0	1	0
			4487	2880	757	826	24			
1	B	550	Total	C	N	O	S	0	1	0
			3992	2547	685	738	22			

There are 36 discrepancies between the modelled and reference sequences:

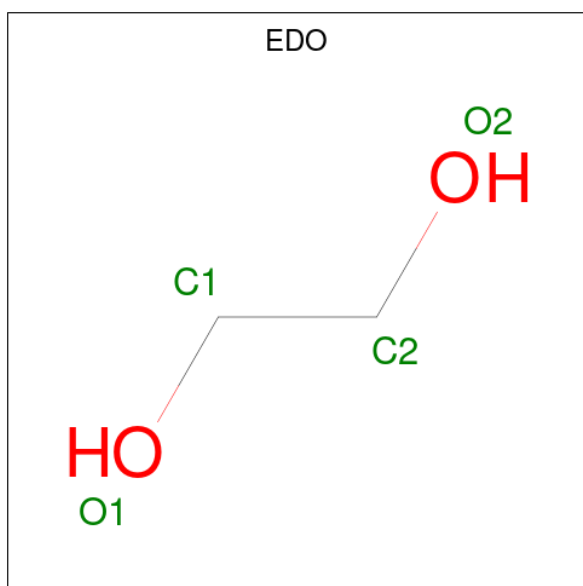
Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	initiating methionine	UNP P35606
A	-16	HIS	-	expression tag	UNP P35606
A	-15	HIS	-	expression tag	UNP P35606
A	-14	HIS	-	expression tag	UNP P35606
A	-13	HIS	-	expression tag	UNP P35606
A	-12	HIS	-	expression tag	UNP P35606
A	-11	HIS	-	expression tag	UNP P35606
A	-10	SER	-	expression tag	UNP P35606
A	-9	SER	-	expression tag	UNP P35606
A	-8	GLY	-	expression tag	UNP P35606
A	-7	ARG	-	expression tag	UNP P35606
A	-6	GLU	-	expression tag	UNP P35606
A	-5	ASN	-	expression tag	UNP P35606
A	-4	LEU	-	expression tag	UNP P35606
A	-3	TYR	-	expression tag	UNP P35606
A	-2	PHE	-	expression tag	UNP P35606
A	-1	GLN	-	expression tag	UNP P35606
A	0	GLY	-	expression tag	UNP P35606
B	-17	MET	-	initiating methionine	UNP P35606
B	-16	HIS	-	expression tag	UNP P35606
B	-15	HIS	-	expression tag	UNP P35606
B	-14	HIS	-	expression tag	UNP P35606
B	-13	HIS	-	expression tag	UNP P35606
B	-12	HIS	-	expression tag	UNP P35606
B	-11	HIS	-	expression tag	UNP P35606

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	expression tag	UNP P35606
B	-9	SER	-	expression tag	UNP P35606
B	-8	GLY	-	expression tag	UNP P35606
B	-7	ARG	-	expression tag	UNP P35606
B	-6	GLU	-	expression tag	UNP P35606
B	-5	ASN	-	expression tag	UNP P35606
B	-4	LEU	-	expression tag	UNP P35606
B	-3	TYR	-	expression tag	UNP P35606
B	-2	PHE	-	expression tag	UNP P35606
B	-1	GLN	-	expression tag	UNP P35606
B	0	GLY	-	expression tag	UNP P35606

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total	O	0	0
			48	48		
3	B	32	Total	O	0	0
			32	32		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.07Å 134.06Å 144.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.12 – 2.40 49.12 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.12-2.40) 98.7 (49.12-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.39Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.219 , 0.247 0.231 , 0.264	Depositor DCC
R_{free} test set	2516 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	52.1	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8563	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: EDO

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	0.52	0/4600	0.73	0/6269
1	B	0.49	0/4084	0.73	0/5578
All	All	0.51	0/8684	0.73	0/11847

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	4487	0	4173	29	0
1	B	3992	0	3467	21	0
2	A	4	0	6	1	0
3	A	48	0	0	0	0
3	B	32	0	0	0	0
All	All	8563	0	7646	50	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 (50) 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:B:9:ARG:HD2	1:B:295:ILE:HD12	1.77	0.67
1:A:535:ILE:HG21	1:A:573:LEU:HD21	1.82	0.60
1:A:358:ILE:HG12	1:A:369:VAL:HG22	1.82	0.59
1:B:114:ILE:HG12	1:B:132:TRP:HZ3	1.66	0.58
1:B:274:VAL:HG22	1:B:290:TYR:CE1	2.41	0.56
1:B:572:ARG:HG2	1:B:584:SER:HB2	1.87	0.55
1:B:31:LEU:HD21	1:B:63:ALA:HB3	1.89	0.55
1:A:31:LEU:HD21	1:A:63:ALA:HB3	1.91	0.53
1:A:17:ARG:HG3	1:A:291:ASP:HA	1.91	0.52
1:A:276:CYS:HB3	2:A:601:EDO:H22	1.92	0.52
1:A:41:VAL:HB	1:A:51:LYS:HB2	1.92	0.52
1:A:271:MET:HB2	1:A:290:TYR:CD2	2.44	0.52
1:A:141:HIS:CE1	1:A:168:LYS:HD2	2.45	0.51
1:B:17:ARG:HG3	1:B:291:ASP:HA	1.92	0.51
1:A:230:GLN:HG2	1:A:250:ASP:HB3	1.91	0.51
1:B:141:HIS:CE1	1:B:168:LYS:HD2	2.46	0.51
1:B:271:MET:HE2	1:B:295:ILE:HG12	1.93	0.50
1:A:324:GLN:HB2	1:A:348:MET:SD	2.51	0.50
1:A:561:MET:HE3	1:A:583:ILE:HD13	1.94	0.50
1:B:41:VAL:HB	1:B:51:LYS:HB2	1.94	0.50
1:B:158:PHE:CE2	1:B:170:TRP:HB2	2.48	0.49
1:A:58:LEU:HD12	1:A:78:ASP:HB3	1.96	0.48
1:A:274:VAL:HG22	1:A:290:TYR:CE2	2.49	0.48
1:A:557:LEU:HD13	1:A:561:MET:HB2	1.96	0.47
1:A:409:ASN:HA	1:A:430:ALA:O	2.15	0.47
1:A:272:GLU:HB2	1:A:291:ASP:HB2	1.97	0.47
1:A:19:LYS:HA	1:A:19:LYS:HE2	1.97	0.46
1:A:312:GLY:HA3	1:A:567:ILE:HD11	1.99	0.45
1:A:37:GLY:HA2	1:A:60:VAL:HG23	2.00	0.44
1:A:524:VAL:HG13	1:A:536:TYR:HB2	1.99	0.44
1:B:17:ARG:HB3	1:B:35:TYR:HB2	1.98	0.44
1:B:557:LEU:HD13	1:B:561:MET:HB2	1.99	0.44
1:A:55:VAL:HG13	1:A:82:ARG:HD2	1.98	0.44
1:B:374:GLU:HA	1:B:390:SER:HA	1.99	0.44
1:A:17:ARG:HB3	1:A:35:TYR:HB2	2.00	0.43
1:B:19:LYS:HA	1:B:19:LYS:HE2	1.99	0.43
1:B:211:VAL:HG21	1:B:246:THR:HG21	2.00	0.43
1:B:112:PRO:HA	1:B:128:TRP:CZ2	2.53	0.43
1:B:228:HIS:CE1	1:B:254[A]:ARG:HG3	2.54	0.43
1:B:357:THR:OG1	1:B:370:CYS:HB2	2.19	0.43
1:A:572:ARG:HG2	1:A:584:SER:HB2	2.00	0.43
1:A:112:PRO:HA	1:A:128:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:LEU:HD11	1:B:515:GLU:HG2	2.00	0.42
1:A:357:THR:OG1	1:A:370:CYS:HB2	2.20	0.42
1:A:211:VAL:HG21	1:A:246:THR:HG21	2.01	0.42
1:B:241:LEU:HB3	1:B:243:ILE:HG12	2.03	0.41
1:A:241:LEU:HB3	1:A:243:ILE:HG12	2.02	0.41
1:A:30:MET:HB2	1:A:296:ILE:HD11	2.03	0.40
1:A:393:GLU:HB2	1:A:406:ARG:HE	1.87	0.40
1:B:70:ASN:HA	1:B:86:TYR:CZ	2.56	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	583/615 (95%)	553 (95%)	27 (5%)	3 (0%)	29	41
1	B	531/615 (86%)	501 (94%)	28 (5%)	2 (0%)	34	48
All	All	1114/1230 (91%)	1054 (95%)	55 (5%)	5 (0%)	34	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	401	SER
1	A	511	GLU
1	B	272	GLU
1	A	372	ASP
1	B	372	ASP

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	444/537 (83%)	437 (98%)	7 (2%)	62	79
1	B	354/537 (66%)	345 (98%)	9 (2%)	47	67
All	All	798/1074 (74%)	782 (98%)	16 (2%)	55	74

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	31	LEU
1	A	217	GLN
1	A	230	GLN
1	A	295	ILE
1	A	316	TRP
1	A	451	ASP
1	B	11	LEU
1	B	31	LEU
1	B	136	GLN
1	B	217	GLN
1	B	271	MET
1	B	316	TRP
1	B	401	SER
1	B	542	ARG
1	B	572	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	359	GLN
1	A	571	ASN

5.3.3 RNA ⓘ

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](#)

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	EDO	A	601	-	3,3,3	0.89	0	2,2,2	0.15	0

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	EDO	A	601	-	-	0/1/1/1	-

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	EDO	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 ⓘ

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	586/615 (95%)	-0.25	1 (0%) 95 94	31, 52, 79, 140	0
1	B	550/615 (89%)	0.16	29 (5%) 26 25	33, 66, 134, 160	0
All	All	1136/1230 (92%)	-0.05	30 (2%) 56 54	31, 58, 111, 160	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	437	PHE	4.7
1	B	458	ILE	4.4
1	B	493	GLU	4.4
1	B	404	ALA	4.3
1	B	452	TRP	4.1
1	B	346	LYS	4.0
1	B	491	LEU	3.8
1	B	492	SER	3.8
1	B	395	ALA	3.5
1	B	490	TYR	3.5
1	B	438	LEU	3.4
1	B	495	VAL	3.4
1	B	375	TYR	3.3
1	B	412	VAL	3.0
1	B	487	ILE	2.9
1	B	461	ILE	2.9
1	B	450	TYR	2.9
1	B	449	PHE	2.8
1	B	488	LEU	2.7
1	A	504	GLY	2.6
1	B	325	ALA	2.5
1	B	515	GLU	2.3
1	B	403	TYR	2.3
1	B	514	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	457	LEU	2.2
1	B	434	TYR	2.2
1	B	382	ALA	2.1
1	B	448	ALA	2.1
1	B	470	TRP	2.0
1	B	405	ILE	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
2	EDO	A	601	4/4	0.93	0.33	52,57,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.