



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:34 PM GMT

PDB ID : 4CNU
Title : CRYSTAL STRUCTURE OF WT HUMAN CRMP-4 from lattice translocation
Authors : Ponnusamy, R.; Lebedev, A.; Lohkamp, B.
Deposited on : 2014-01-24
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

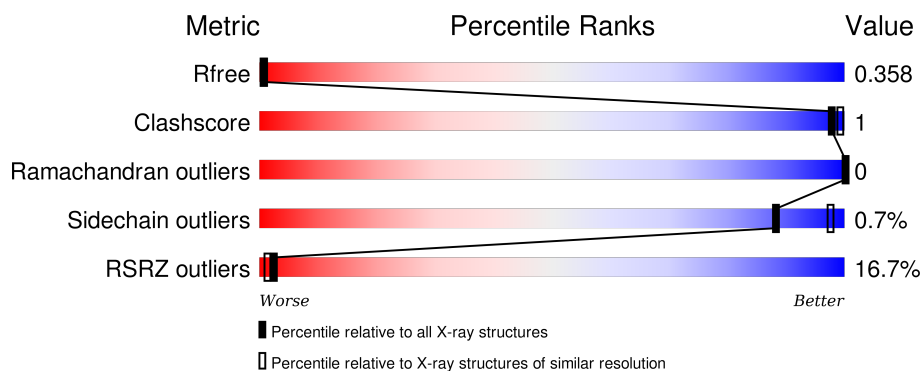
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	570	<div> <div>8%</div> <div>83%</div> <div>14%</div> </div>
1	B	570	<div> <div>21%</div> <div>82%</div> <div>14%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7512 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 DIHYDROPYRIMIDINASE-LIKE 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	0	0
			3756	2373	635	725	23			
1	B	488	Total	C	N	O	S	0	0	0
			3756	2373	635	725	23			

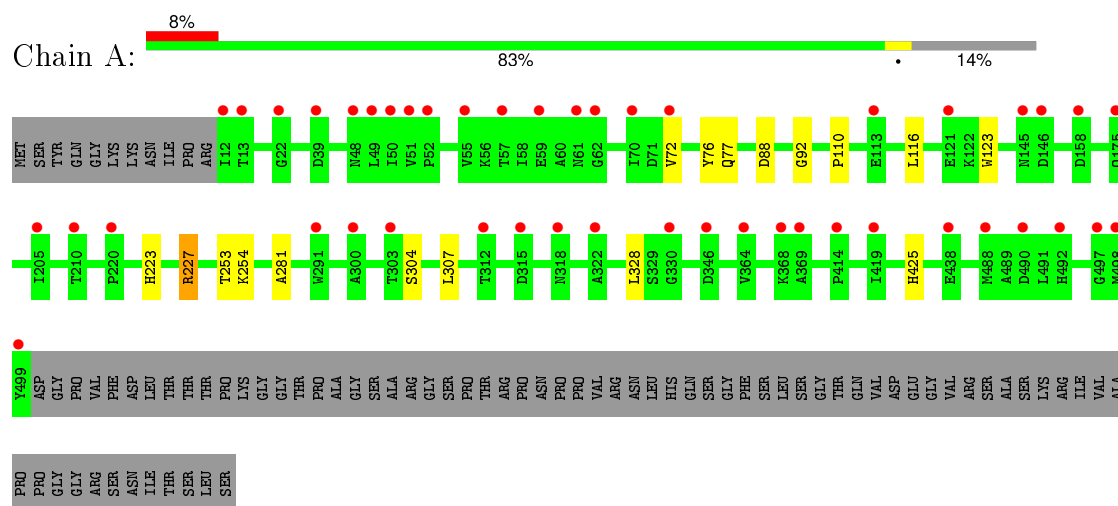
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	TYR	PHE	CONFLICT	UNP Q14195
B	76	TYR	PHE	CONFLICT	UNP Q14195

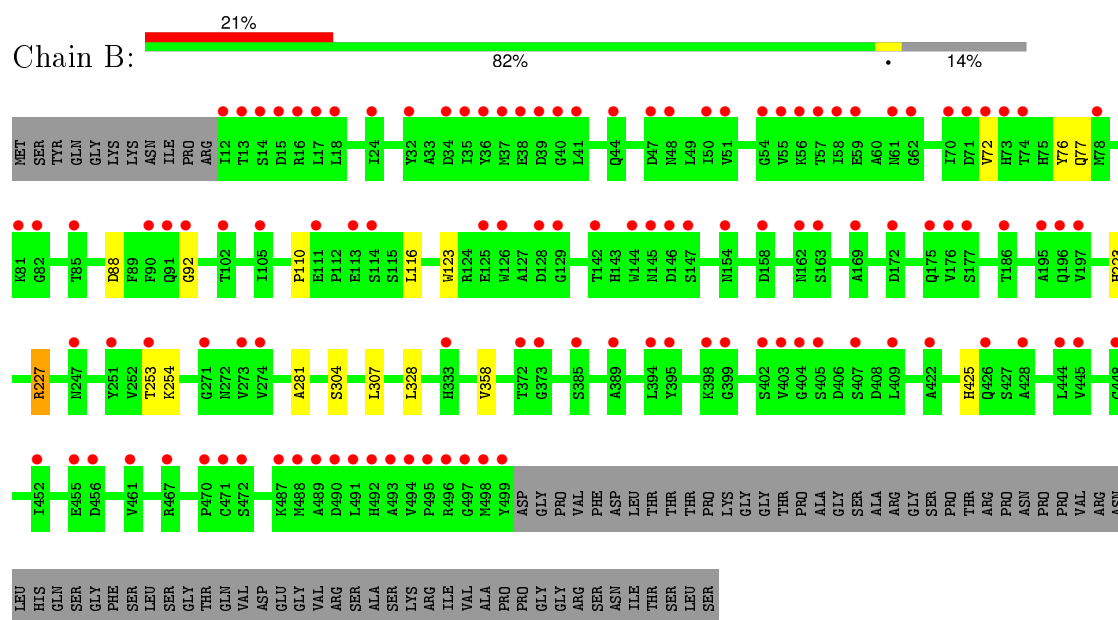
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: DIHYDROPYRIMIDINASE-LIKE 3



• Molecule 1: DIHYDROPYRIMIDINASE-LIKE 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	96.24Å 108.22Å 118.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.43 – 2.80 39.40 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.0 (39.43-2.80) 93.1 (39.40-2.80)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.255 , 0.295 0.322 , 0.358	Depositor DCC
R_{free} test set	1465 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.570	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 25.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	13 of 28919 reflections (0.045%)	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7512	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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 68.08 % 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 4.6339e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.31	0/3835	0.50	0/5201
1	B	0.31	0/3835	0.50	0/5201
All	All	0.31	0/7670	0.50	0/10402

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	3756	0	3700	8	0
1	B	3756	0	3700	9	0
All	All	7512	0	7400	17	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 (17) 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:76:TYR:CE1	1:B:92:GLY:HA3	2.43	0.53
1:A:76:TYR:CE1	1:A:92:GLY:HA3	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:HIS:O	1:A:227:ARG:HG3	2.11	0.51
1:B:223:HIS:O	1:B:227:ARG:HG3	2.13	0.49
1:B:77:GLN:OE1	1:B:88:ASP:HB2	2.14	0.48
1:A:77:GLN:OE1	1:A:88:ASP:HB2	2.14	0.47
1:A:253:THR:C	1:A:254:LYS:HG2	2.36	0.45
1:B:110:PRO:HG3	1:B:116:LEU:HG	1.99	0.45
1:A:110:PRO:HG3	1:A:116:LEU:HG	1.99	0.45
1:B:281:ALA:HB1	1:B:307:LEU:HD11	1.98	0.45
1:A:281:ALA:HB1	1:A:307:LEU:HD11	1.99	0.44
1:B:253:THR:C	1:B:254:LYS:HG2	2.39	0.42
1:A:72:VAL:HG11	1:A:328:LEU:HD12	2.02	0.42
1:B:72:VAL:HG11	1:B:328:LEU:HD12	2.02	0.41
1:B:76:TYR:OH	1:B:358:VAL:HG21	2.20	0.41
1:A:254:LYS:CE	1:A:304:SER:O	2.68	0.41
1:B:254:LYS:CE	1:B:304:SER:O	2.69	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	486/570 (85%)	461 (95%)	25 (5%)	0	100	100
1	B	486/570 (85%)	463 (95%)	23 (5%)	0	100	100
All	All	972/1140 (85%)	924 (95%)	48 (5%)	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	404/471 (86%)	401 (99%)	3 (1%)	88	97
1	B	404/471 (86%)	401 (99%)	3 (1%)	88	97
All	All	808/942 (86%)	802 (99%)	6 (1%)	88	97

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	TRP
1	A	227	ARG
1	A	425	HIS
1	B	123	TRP
1	B	227	ARG
1	B	425	HIS

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

Mol	Chain	Res	Type
1	A	29	GLN
1	B	29	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	488/570 (85%)	0.69	46 (9%) 11 5	26, 42, 66, 108	0
1	B	488/570 (85%)	1.35	117 (23%) 1 1	25, 46, 75, 99	0
All	All	976/1140 (85%)	1.02	163 (16%) 2 1	25, 44, 72, 108	0

All (163) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	VAL	9.9
1	B	55	VAL	9.4
1	B	489	ALA	7.7
1	B	51	VAL	7.4
1	B	48	ASN	6.6
1	A	49	LEU	6.6
1	B	492	HIS	6.0
1	B	61	ASN	5.6
1	B	495	PRO	5.3
1	B	125	GLU	5.0
1	B	488	MET	5.0
1	B	40	GLY	4.9
1	B	491	LEU	4.7
1	A	55	VAL	4.6
1	A	318	ASN	4.6
1	B	247	ASN	4.6
1	B	452	ILE	4.5
1	B	54	GLY	4.4
1	B	62	GLY	4.3
1	B	72	VAL	4.2
1	B	146	ASP	4.2
1	A	146	ASP	4.1
1	B	158	ASP	4.1
1	B	57	THR	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	32	TYR	3.9
1	A	490	ASP	3.8
1	B	490	ASP	3.8
1	A	61	ASN	3.7
1	B	13	THR	3.7
1	A	13	THR	3.7
1	A	364	VAL	3.7
1	A	48	ASN	3.6
1	B	407	SER	3.6
1	A	312	THR	3.6
1	B	70	ILE	3.5
1	B	15	ASP	3.5
1	B	92	GLY	3.5
1	B	82	GLY	3.4
1	A	158	ASP	3.4
1	A	50	ILE	3.4
1	B	176	VAL	3.4
1	A	492	HIS	3.4
1	B	445	VAL	3.4
1	B	186	THR	3.4
1	B	129	GLY	3.4
1	A	12	ILE	3.4
1	B	12	ILE	3.4
1	B	35	ILE	3.4
1	B	18	LEU	3.4
1	B	455	GLU	3.3
1	B	177	SER	3.3
1	B	59	GLU	3.3
1	B	385	SER	3.3
1	A	300	ALA	3.3
1	B	39	ASP	3.3
1	B	487	LYS	3.3
1	B	498	MET	3.2
1	B	497	GLY	3.2
1	B	74	THR	3.2
1	A	488	MET	3.1
1	B	496	ARG	3.0
1	A	59	GLU	3.0
1	A	22	GLY	3.0
1	B	91	GLN	3.0
1	B	90	PHE	2.9
1	B	147	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	36	TYR	2.9
1	B	172	ASP	2.9
1	B	44	GLN	2.8
1	A	210	THR	2.8
1	A	62	GLY	2.8
1	B	38	GLU	2.8
1	A	39	ASP	2.8
1	B	472	SER	2.8
1	B	493	ALA	2.7
1	A	498	MET	2.7
1	B	499	TYR	2.7
1	B	373	GLY	2.7
1	B	114	SER	2.7
1	A	497	GLY	2.7
1	B	253	THR	2.7
1	B	461	VAL	2.7
1	A	113	GLU	2.6
1	A	175	GLN	2.6
1	B	145	ASN	2.6
1	B	470	PRO	2.6
1	A	57	THR	2.6
1	A	72	VAL	2.6
1	B	426	GLN	2.6
1	A	315	ASP	2.6
1	B	404	GLY	2.5
1	B	409	LEU	2.5
1	B	467	ARG	2.5
1	B	394	LEU	2.5
1	B	399	GLY	2.5
1	B	471	CYS	2.5
1	B	105	ILE	2.5
1	B	56	LYS	2.4
1	B	154	ASN	2.4
1	B	372	THR	2.4
1	B	128	ASP	2.4
1	A	303	THR	2.4
1	B	448	CYS	2.4
1	B	422	ALA	2.4
1	A	368	LYS	2.4
1	B	494	VAL	2.4
1	B	16	ARG	2.4
1	B	17	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	37	MET	2.4
1	A	322	ALA	2.4
1	A	70	ILE	2.4
1	B	24	ILE	2.4
1	A	499	TYR	2.4
1	B	389	ALA	2.3
1	A	346	ASP	2.3
1	B	456	ASP	2.3
1	B	175	GLN	2.3
1	B	398	LYS	2.3
1	B	113	GLU	2.3
1	B	162	ASN	2.3
1	B	169	ALA	2.3
1	A	52	PRO	2.3
1	B	34	ASP	2.3
1	A	369	ALA	2.3
1	B	405	SER	2.3
1	A	438	GLU	2.3
1	B	403	VAL	2.3
1	B	81	LYS	2.3
1	B	50	ILE	2.2
1	B	195	ALA	2.2
1	B	395	TYR	2.2
1	A	330	GLY	2.2
1	B	71	ASP	2.2
1	B	428	ALA	2.2
1	B	163	SER	2.2
1	B	197	VAL	2.2
1	B	273	VAL	2.2
1	B	126	TRP	2.2
1	B	144	TRP	2.2
1	B	402	SER	2.2
1	B	102	THR	2.2
1	A	291	TRP	2.2
1	B	444	LEU	2.1
1	A	205	ILE	2.1
1	B	251	TYR	2.1
1	B	47	ASP	2.1
1	A	121	GLU	2.1
1	A	414	PRO	2.1
1	B	274	VAL	2.1
1	B	271	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	78	MET	2.1
1	B	85	THR	2.1
1	A	220	PRO	2.1
1	B	41	LEU	2.1
1	B	111	GLU	2.1
1	B	333	HIS	2.0
1	B	58	ILE	2.0
1	A	145	ASN	2.0
1	B	14	SER	2.0
1	B	73	HIS	2.0
1	B	142	THR	2.0
1	A	419	ILE	2.0
1	B	196	GLN	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 carbohydrates 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.