



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5Q4N / pdb_00005q4n
Title : PanDDA analysis group deposition – Crystal Structure of DCLRE1A after initial refinement with no ligand modelled (structure 88)
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Deposited on : 2017-05-25
Resolution : 1.71 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
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

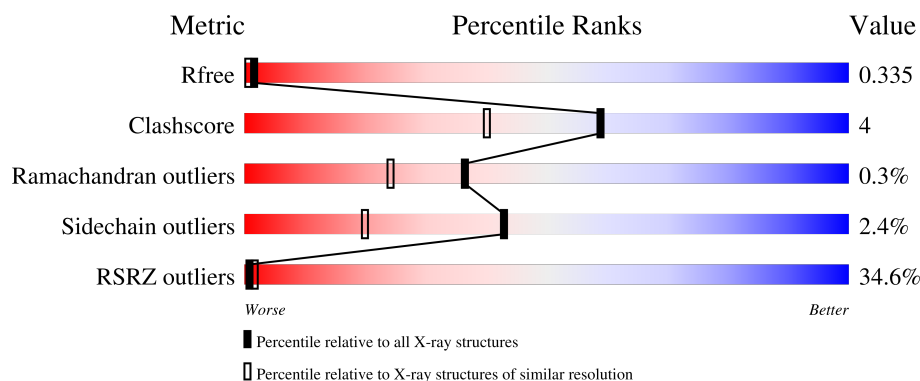
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	343	34% 88% 11% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3044 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 DCLRE1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	341	2726	1765	450	491	20	0	4	0

- Molecule 2 is MALONATE ION (CCD ID: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	7	3	4	0	0

- Molecule 3 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ni	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	310	Total 310	O 310	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.93Å 57.94Å 114.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.69 – 1.71 23.69 – 1.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (23.69-1.71) 99.9 (23.69-1.71)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 1.71Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.269 , 0.324 0.281 , 0.335	Depositor DCC
R_{free} test set	1865 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.3	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	0.93	EDS
Total number of atoms	3044	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MLI, NI

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	0.98	0/2802	1.05	3/3809 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	860	ASN	N-CA-C	6.68	118.25	110.97
1	A	848	PHE	CB-CA-C	5.15	117.02	110.22
1	A	830	GLN	N-CA-C	5.04	116.51	108.79

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	2726	0	2676	24	0
2	A	7	0	2	0	0
3	A	1	0	0	0	0
4	A	310	0	0	14	0
All	All	3044	0	2678	24	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 4.

All (24) 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:A:909:GLN:NE2	4:A:1201:HOH:O	2.09	0.86
1:A:1032:ARG:NH2	4:A:1209:HOH:O	2.31	0.64
1:A:896:SER:O	4:A:1203:HOH:O	2.15	0.64
1:A:815:ASP:OD1	4:A:1202:HOH:O	2.15	0.62
1:A:1032:ARG:NE	4:A:1214:HOH:O	2.40	0.54
1:A:1035:LYS:NZ	4:A:1213:HOH:O	2.40	0.54
1:A:906:LYS:NZ	4:A:1212:HOH:O	2.39	0.53
1:A:743:LYS:HG3	1:A:744:HIS:CE1	2.45	0.51
1:A:973:ILE:HG23	4:A:1443:HOH:O	2.11	0.50
1:A:1032:ARG:CZ	4:A:1214:HOH:O	2.59	0.49
1:A:874:VAL:HG11	1:A:890:ILE:HG21	1.93	0.48
1:A:717:PHE:CE1	1:A:728:TYR:HB3	2.48	0.48
1:A:742:SER:O	1:A:765:VAL:HG22	2.13	0.47
1:A:718:GLN:HG2	4:A:1285:HOH:O	2.18	0.43
1:A:720:GLY:N	4:A:1227:HOH:O	2.51	0.43
1:A:860:ASN:O	1:A:861:THR:C	2.62	0.43
1:A:718:GLN:HB3	1:A:739:ALA:O	2.19	0.42
1:A:856:ARG:HE	1:A:856:ARG:HA	1.85	0.42
1:A:889:ALA:O	4:A:1204:HOH:O	2.22	0.42
1:A:875:VAL:HG23	1:A:956:ILE:HG23	2.02	0.42
1:A:851:GLN:NE2	4:A:1231:HOH:O	2.52	0.42
1:A:760:LYS:O	1:A:764[A]:HIS:HA	2.20	0.41
1:A:898:VAL:HG11	1:A:931:LEU:HD21	2.02	0.41
1:A:984:ILE:HD12	4:A:1300:HOH:O	2.21	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	343/343 (100%)	326 (95%)	16 (5%)	1 (0%)	36 24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	767	GLU

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	299/305 (98%)	292 (98%)	7 (2%)	44 21

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	816	PHE
1	A	830	GLN
1	A	879	TYR
1	A	893	VAL
1	A	928	LEU
1	A	977	ILE
1	A	1029	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	744	HIS
1	A	851	GLN
1	A	909	GLN
1	A	966	HIS
1	A	979	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	MLI	A	1101	3	6,6,6	1.07	0	7,7,7	0.98	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	MLI	A	1101	3	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	MLI	C2-C1-C3-O9
2	A	1101	MLI	C2-C1-C3-O8
2	A	1101	MLI	C3-C1-C2-O6
2	A	1101	MLI	C3-C1-C2-O7

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	341/343 (99%)	1.62	118 (34%) 1 2	15, 37, 104, 174	4 (1%)

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	973	ILE	8.1
1	A	984	ILE	7.6
1	A	970	PHE	7.1
1	A	765	VAL	7.0
1	A	939	PHE	7.0
1	A	942	LEU	6.7
1	A	986	ILE	6.7
1	A	745	PHE	6.6
1	A	956	ILE	6.6
1	A	977	ILE	6.5
1	A	971	THR	6.4
1	A	982	GLY	6.2
1	A	744	HIS	6.2
1	A	980	THR	6.0
1	A	974	ALA	5.9
1	A	937	ILE	5.7
1	A	769	TYR	5.6
1	A	978	PRO	5.6
1	A	976	VAL	5.4
1	A	966	HIS	5.4
1	A	946	LEU	5.4
1	A	967	SER	5.3
1	A	941	GLY	5.2
1	A	968	ASN	5.2
1	A	985	SER	5.0
1	A	987	TYR	5.0
1	A	951	GLY	4.9

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Mol	Chain	Res	Type	RSRZ
1	A	954	ASN	4.8
1	A	747	PHE	4.8
1	A	781	VAL	4.8
1	A	972	ARG	4.7
1	A	711	GLY	4.7
1	A	981	LYS	4.6
1	A	964	TRP	4.5
1	A	780	ILE	4.5
1	A	950	GLY	4.4
1	A	768	GLN	4.2
1	A	928	LEU	4.1
1	A	953	TYR	4.1
1	A	746	THR	4.1
1	A	983	ASN	4.1
1	A	866	VAL	3.8
1	A	965	THR	3.8
1	A	724	GLY	3.8
1	A	949	CYS	3.8
1	A	770	ILE	3.7
1	A	963	GLY	3.7
1	A	722	VAL	3.7
1	A	871	HIS	3.7
1	A	742	SER	3.7
1	A	940	LYS	3.5
1	A	862	ALA	3.5
1	A	764[A]	HIS	3.5
1	A	750	TYR	3.5
1	A	938	ASN	3.5
1	A	868	LEU	3.4
1	A	873	LEU	3.4
1	A	741	LEU	3.4
1	A	700	THR	3.3
1	A	763	LEU	3.3
1	A	784	VAL	3.3
1	A	957	LEU	3.2
1	A	879	TYR	3.2
1	A	988	GLY	3.2
1	A	743	LYS	3.2
1	A	943	GLN	3.1
1	A	969	LYS	3.1
1	A	721	VAL	3.1
1	A	856	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	709	GLY	3.0
1	A	955	GLN	3.0
1	A	975	ASP	3.0
1	A	979	GLN	2.9
1	A	894	LEU	2.9
1	A	861	THR	2.8
1	A	826	LEU	2.8
1	A	927	SER	2.8
1	A	945	HIS	2.8
1	A	766	GLN	2.8
1	A	936	GLN	2.8
1	A	867	THR	2.8
1	A	767	GLU	2.7
1	A	863	PHE	2.7
1	A	748	PRO	2.7
1	A	935	MET	2.7
1	A	779	CYS	2.7
1	A	925	CYS	2.7
1	A	932	LEU	2.7
1	A	740	GLY	2.7
1	A	865	ALA	2.7
1	A	870	PRO	2.6
1	A	848	PHE	2.6
1	A	782	ASN	2.6
1	A	926	SER	2.6
1	A	962	THR	2.5
1	A	874	VAL	2.5
1	A	893	VAL	2.5
1	A	989	ILE	2.4
1	A	952	LYS	2.3
1	A	869	ASN	2.3
1	A	749	VAL	2.3
1	A	961	PRO	2.3
1	A	739	ALA	2.2
1	A	948	LYS	2.2
1	A	857	PHE	2.2
1	A	726	THR	2.2
1	A	708	PRO	2.2
1	A	710	THR	2.2
1	A	875	VAL	2.1
1	A	771	HIS	2.1
1	A	895	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	759	LEU	2.1
1	A	717	PHE	2.1
1	A	723	GLU	2.1
1	A	944	SER	2.1
1	A	872	ALA	2.0
1	A	725	CYS	2.0
1	A	878	THR	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	MLI	A	1101	7/7	0.94	0.07	23,31,33,34	0
3	NI	A	1102	1/1	0.98	0.04	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.