



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:32 PM UTC

PDB ID : 5Q7C / pdb_00005q7c
Title : PanDDA analysis group deposition – Crystal Structure of DCLRE1A after initial refinement with no ligand modelled (structure 185)
Authors : Newman, J.A.; Aitkenhead, H.; Lee, S.Y.; Kupinska, K.; Burgess-Brown, N.; Tallon, R.; Krojer, T.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.; Bountra, C.; Gileadi, O.
Deposited on : 2017-05-25
Resolution : 1.20 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

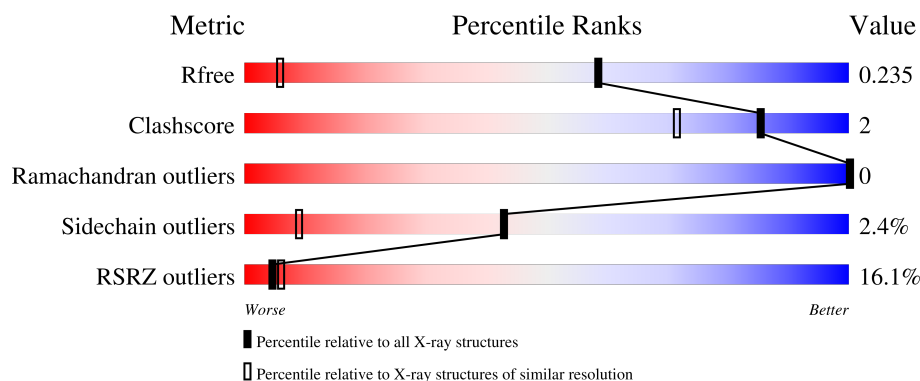
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.20 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	1216 (1.20-1.20)
Clashscore	190562	1265 (1.20-1.20)
Ramachandran outliers	187476	1226 (1.20-1.20)
Sidechain outliers	187428	1226 (1.20-1.20)
RSRZ outliers	180081	1214 (1.20-1.20)

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	

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
2	MLI	A	1101	-	X	-	-

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

- Molecule 1: DCLRE1A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.95Å 57.02Å 115.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	115.23 – 1.20 57.62 – 1.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (115.23-1.20) 99.1 (57.62-1.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 1.20Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.211 , 0.232 (Not available) , 0.235	Depositor DCC
R_{free} test set	5292 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	13.7	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3044	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% 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	1.78	34/2802 (1.2%)	1.41	14/3809 (0.4%)

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	989	ILE	CA-C	11.15	1.60	1.52
1	A	707	ILE	CA-C	8.95	1.60	1.53
1	A	891	ALA	C-O	7.39	1.32	1.24
1	A	1006	TRP	C-O	7.10	1.32	1.24
1	A	828	ALA	N-CA	6.87	1.54	1.46
1	A	1032	ARG	N-CA	6.86	1.54	1.46
1	A	961	PRO	N-CA	-6.81	1.39	1.47
1	A	777	THR	N-CA	6.73	1.54	1.46
1	A	1023	SER	C-O	6.40	1.31	1.24
1	A	892	ASP	C-O	6.37	1.31	1.24
1	A	888	LEU	C-O	6.21	1.31	1.24
1	A	860	ASN	N-CA	6.14	1.53	1.46
1	A	988	GLY	C-N	6.00	1.37	1.33
1	A	808	THR	N-CA	5.95	1.53	1.46
1	A	921	THR	N-CA	5.94	1.53	1.46
1	A	930	HIS	N-CA	5.89	1.53	1.46
1	A	1025	SER	C-O	5.88	1.31	1.24
1	A	956	ILE	N-CA	5.86	1.53	1.46
1	A	892	ASP	N-CA	5.79	1.53	1.46
1	A	1026	THR	C-O	5.74	1.30	1.24
1	A	890	ILE	C-O	5.55	1.30	1.24
1	A	1020	THR	C-O	5.45	1.31	1.23
1	A	990	PRO	CA-C	5.33	1.57	1.53
1	A	879	TYR	N-CA	5.30	1.51	1.46
1	A	811	LEU	N-CA	5.27	1.52	1.46
1	A	927	SER	C-O	-5.25	1.17	1.24

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	761	ASN	N-CA	5.22	1.52	1.46
1	A	760	LYS	N-CA	5.18	1.52	1.46
1	A	828	ALA	CA-C	5.17	1.59	1.52
1	A	835	LEU	N-CA	5.12	1.52	1.46
1	A	859	ILE	N-CA	5.09	1.52	1.46
1	A	824	ARG	N-CA	5.08	1.52	1.46
1	A	865	ALA	N-CA	-5.05	1.40	1.46
1	A	889	ALA	C-O	5.01	1.30	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	829	ASP	N-CA-C	9.78	124.61	112.87
1	A	961	PRO	CB-CA-C	8.06	118.73	111.40
1	A	961	PRO	N-CA-C	-7.71	97.56	110.21
1	A	829	ASP	CB-CA-C	-6.19	96.66	110.21
1	A	766	GLN	CA-C-N	6.12	128.77	120.38
1	A	766	GLN	C-N-CA	6.12	128.77	120.38
1	A	898	VAL	CA-C-N	5.51	126.92	121.57
1	A	898	VAL	C-N-CA	5.51	126.92	121.57
1	A	911	LEU	N-CA-C	5.30	119.44	112.92
1	A	900	MET	CA-C-O	-5.27	115.80	121.38
1	A	828	ALA	CA-C-N	5.19	130.65	120.99
1	A	828	ALA	C-N-CA	5.19	130.65	120.99
1	A	887	PHE	N-CA-C	-5.19	105.69	112.23
1	A	1004	VAL	O-C-N	5.17	127.15	121.83

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	12	1
2	A	7	0	2	0	0
3	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	310	0	0	8	0
All	All	3044	0	2678	12	1

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 2.

All (12) 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:936:GLN:OE1	4:A:1201:HOH:O	2.05	0.74
1:A:909:GLN:NE2	4:A:1202:HOH:O	2.31	0.63
1:A:973:ILE:HG22	4:A:1445:HOH:O	2.04	0.57
1:A:1018:VAL:HG22	4:A:1239:HOH:O	2.03	0.56
1:A:935:MET:HE2	1:A:935:MET:HB3	1.80	0.46
1:A:869:ASN:ND2	1:A:871:HIS:H	2.14	0.46
1:A:935:MET:HG3	4:A:1430:HOH:O	2.16	0.45
1:A:1032:ARG:CZ	4:A:1220:HOH:O	2.64	0.45
1:A:1032:ARG:NH2	4:A:1209:HOH:O	2.50	0.44
1:A:830:GLN:HB2	4:A:1414:HOH:O	2.17	0.43
1:A:875:VAL:O	1:A:958:ALA:HA	2.21	0.40
1:A:743:LYS:CB	1:A:764[B]:HIS:O	2.70	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:829:ASP:OD2	1:A:966:HIS:N[3_545]	2.18	0.02

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%)	337 (98%)	6 (2%)	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	299/305 (98%)	292 (98%)	7 (2%)	44	9

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	856	ARG
1	A	879	TYR
1	A	928	LEU
1	A	960	ARG
1	A	1029	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	768	GLN
1	A	851	GLN
1	A	860	ASN
1	A	869	ASN
1	A	966	HIS
1	A	968	ASN
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 [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	1.81	3 (50%)	7,7,7	0.77	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	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1101	MLI	C1-C2	2.48	1.54	1.51
2	A	1101	MLI	C1-C3	2.42	1.54	1.51
2	A	1101	MLI	O7-C2	-2.40	1.22	1.30

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	C3-C1-C2-O6
2	A	1101	MLI	C2-C1-C3-O8
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%)	0.91	55 (16%) 4 6	7, 19, 37, 72	4 (1%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	970	PHE	5.7
1	A	879	TYR	5.1
1	A	703	PHE	4.4
1	A	964	TRP	4.2
1	A	971	THR	4.1
1	A	721	VAL	3.9
1	A	700	THR	3.8
1	A	925	CYS	3.6
1	A	966	HIS	3.5
1	A	969	LYS	3.4
1	A	968	ASN	3.4
1	A	928	LEU	3.3
1	A	740	GLY	3.3
1	A	829	ASP	3.3
1	A	950	GLY	3.2
1	A	830	GLN	3.2
1	A	963	GLY	3.2
1	A	965	THR	3.1
1	A	916	ILE	3.0
1	A	977	ILE	3.0
1	A	708	PRO	3.0
1	A	769	TYR	3.0
1	A	832	VAL	3.0
1	A	742	SER	2.9
1	A	747	PHE	2.8
1	A	967	SER	2.8
1	A	720	GLY	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	709	GLY	2.8
1	A	710	THR	2.7
1	A	746	THR	2.7
1	A	741	LEU	2.7
1	A	724	GLY	2.7
1	A	962	THR	2.7
1	A	745	PHE	2.7
1	A	764[A]	HIS	2.6
1	A	826	LEU	2.5
1	A	957	LEU	2.4
1	A	961	PRO	2.4
1	A	973	ILE	2.3
1	A	704	TYR	2.3
1	A	936	GLN	2.2
1	A	913	ILE	2.2
1	A	726	THR	2.2
1	A	935	MET	2.2
1	A	972	ARG	2.1
1	A	979	GLN	2.1
1	A	828	ALA	2.1
1	A	960	ARG	2.1
1	A	951	GLY	2.1
1	A	701	CYS	2.1
1	A	744	HIS	2.1
1	A	974	ALA	2.1
1	A	926	SER	2.0
1	A	707	ILE	2.0
1	A	917	ASN	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	0.96	0.07	14,17,22,29	0
3	NI	A	1102	1/1	1.00	0.05	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.