



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:56 AM UTC

PDB ID : 7EIP / pdb_00007eip
Title : Crystal structure of ligand-free chondroitin ABC lyase I
Authors : Takashima, M.; Miyanaga, A.; Eguchi, T.
Deposited on : 2021-03-31
Resolution : 1.88 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

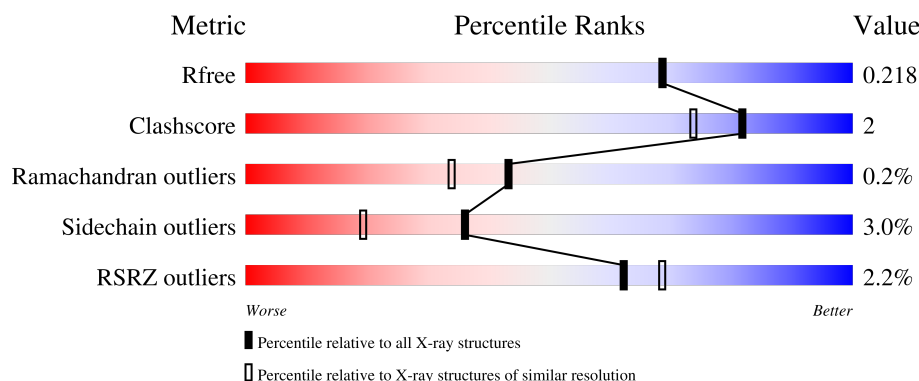
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

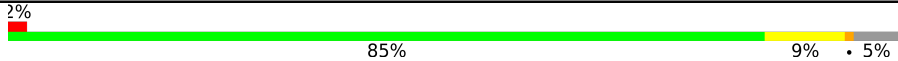
The reported resolution of this entry is 1.88 Å.

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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	1021	

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	ACT	A	1102	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8530 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 Chondroitin sulfate ABC endolyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	966	Total	C	N	O	S	0	3	0
			7734	4922	1319	1472	21			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	787	Total 787	O 787	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	48.94Å 94.24Å 229.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.47 – 1.88 43.47 – 1.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.47-1.88) 100.0 (43.47-1.88)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.169 , 0.213 0.179 , 0.218	Depositor DCC
R_{free} test set	4349 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	21.3	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8530	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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: ACT, 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	1.15	14/7932 (0.2%)	1.33	8/10757 (0.1%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	619	PRO	C-O	-6.66	1.15	1.23
1	A	798	TRP	C-O	6.30	1.31	1.24
1	A	673	VAL	C-O	6.26	1.30	1.24
1	A	328	GLN	C-O	6.13	1.31	1.24
1	A	407	ALA	C-O	-5.91	1.17	1.24
1	A	388	HIS	CE1-NE2	5.68	1.38	1.32
1	A	768	ILE	C-O	5.59	1.29	1.23
1	A	961	ASP	C-O	-5.57	1.17	1.24
1	A	700	HIS	CE1-NE2	5.56	1.38	1.32
1	A	750	PRO	C-O	-5.30	1.17	1.24
1	A	638	LYS	C-O	5.22	1.30	1.23
1	A	1007	SER	C-O	5.20	1.29	1.23
1	A	120	PRO	C-O	-5.18	1.17	1.23
1	A	163	LEU	N-CA	5.09	1.51	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	906	ASP	CA-CB-CG	7.10	119.70	112.60
1	A	252	GLU	CB-CG-CD	6.11	122.99	112.60
1	A	547	ILE	CA-C-O	-6.10	114.39	120.85
1	A	1004	THR	CA-CB-OG1	-5.72	101.02	109.60
1	A	351	VAL	CA-C-O	-5.62	115.21	121.17
1	A	830	THR	N-CA-C	-5.42	107.32	114.04
1	A	338	ASN	N-CA-C	-5.29	105.60	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	TYR	CA-C-O	-5.14	115.82	121.58

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	7734	0	7577	35	0
2	A	8	0	6	3	0
3	A	1	0	0	0	0
4	A	787	0	0	6	1
All	All	8530	0	7583	38	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 (38) 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:637:ASP:OD1	4:A:1201:HOH:O	2.05	0.73
1:A:812:THR:HG22	1:A:836:ASN:HD22	1.54	0.70
1:A:829:ILE:HD13	1:A:835:VAL:HG21	1.81	0.63
2:A:1102:ACT:H1	4:A:1805:HOH:O	1.99	0.62
1:A:812:THR:HG22	1:A:836:ASN:ND2	2.16	0.60
1:A:212:ARG:O	1:A:212:ARG:HD2	2.01	0.60
1:A:58:ASN:HD22	1:A:58:ASN:N	1.98	0.59
2:A:1101:ACT:CH3	4:A:1216:HOH:O	2.53	0.55
1:A:450:SER:HB2	1:A:519:LEU:HD11	1.88	0.54
1:A:954:ILE:HG12	1:A:1016:LYS:HG3	1.91	0.53
1:A:57:LYS:C	1:A:58:ASN:HD22	2.16	0.53
1:A:189:GLY:O	1:A:192:VAL:HG23	2.09	0.52
1:A:257:ILE:HG21	1:A:466:ARG:HD2	1.93	0.50
1:A:734:MET:HA	1:A:758:VAL:O	2.12	0.49
1:A:131:GLU:OE1	1:A:191:LYS:HD3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:GLN:O	1:A:316:ILE:HG22	2.14	0.47
1:A:115:LEU:HA	1:A:209:TYR:O	2.15	0.46
1:A:777:LYS:HB3	1:A:868:PRO:HB3	1.98	0.46
1:A:806:ASN:HB2	4:A:1644:HOH:O	2.14	0.45
1:A:826:GLY:O	1:A:877:VAL:HA	2.17	0.44
1:A:465:LYS:HG2	4:A:1881:HOH:O	2.18	0.44
1:A:798:TRP:O	1:A:819:LEU:HA	2.18	0.43
1:A:282:ASN:OD1	1:A:284:SER:HB2	2.18	0.43
1:A:311:ASP:O	1:A:314:ILE:HG12	2.18	0.43
1:A:735:MET:O	1:A:757:SER:HA	2.19	0.43
1:A:283:ILE:HG21	1:A:358:GLN:HE22	1.84	0.42
1:A:555:LEU:N	1:A:556:PRO:CD	2.83	0.42
1:A:226:ASP:HB3	1:A:383:ALA:HB2	2.01	0.42
1:A:597:ILE:HD12	1:A:597:ILE:HA	1.76	0.42
1:A:264:LEU:HG	1:A:522:LEU:HD13	2.01	0.42
1:A:90:LYS:O	1:A:91:LEU:HD23	2.20	0.41
1:A:451:ARG:HD2	1:A:518:GLN:OE1	2.21	0.41
2:A:1102:ACT:CH3	4:A:1805:HOH:O	2.64	0.41
1:A:664:TYR:CE2	1:A:698:THR:HB	2.55	0.41
1:A:90:LYS:C	1:A:91:LEU:HD23	2.45	0.41
1:A:154:ALA:HB3	1:A:384:LEU:HA	2.03	0.40
1:A:661:TYR:CE1	1:A:703:LEU:HB3	2.56	0.40
1:A:397:TRP:O	1:A:398:TYR:C	2.64	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)
4:A:1511:HOH:O	4:A:1757:HOH:O[3_544]	2.16	0.04

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	959/1021 (94%)	921 (96%)	36 (4%)	2 (0%)	43 34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	950	LYS
1	A	36	LEU

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	850/890 (96%)	825 (97%)	25 (3%)	37 20

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	SER
1	A	58	ASN
1	A	62	THR
1	A	66	LYS
1	A	72	ASN
1	A	76	LEU
1	A	80	LYS
1	A	90	LYS
1	A	119	LYS
1	A	127	ILE
1	A	135	SER
1	A	140	GLN
1	A	188	LEU
1	A	212	ARG
1	A	234	SER
1	A	273	LYS
1	A	371	LYS
1	A	405	SER
1	A	451	ARG

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Mol	Chain	Res	Type
1	A	597	ILE
1	A	599	ASP
1	A	639	MET
1	A	667	HIS
1	A	876	MET
1	A	951	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	58	ASN
1	A	242	ASN
1	A	246	GLN
1	A	296	HIS
1	A	464	GLN
1	A	636	GLN
1	A	648	ASN
1	A	731	GLN
1	A	771	ASN
1	A	836	ASN
1	A	840	GLN
1	A	842	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	ACT	A	1101	-	3,3,3	0.85	0	3,3,3	0.77	0
2	ACT	A	1102	-	3,3,3	0.52	0	3,3,3	1.81	1 (33%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	1102	ACT	O-C-CH3	-2.44	112.52	122.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	ACT	1	0
2	A	1102	ACT	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	966/1021 (94%)	-0.10	21 (2%)	62 68	12, 22, 51, 72	3 (0%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	951	ASP	3.5
1	A	134	ILE	3.2
1	A	133	LEU	3.0
1	A	34	LYS	2.8
1	A	60	ILE	2.8
1	A	137	SER	2.6
1	A	33	PRO	2.5
1	A	392	TYR	2.5
1	A	132	LYS	2.5
1	A	36	LEU	2.4
1	A	164	GLU	2.4
1	A	299	ALA	2.3
1	A	273	LYS	2.3
1	A	188	LEU	2.3
1	A	203	VAL	2.2
1	A	61	LEU	2.2
1	A	332	ASN	2.1
1	A	896	ASN	2.1
1	A	101	LYS	2.1
1	A	135	SER	2.0
1	A	136	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	ACT	A	1102	4/4	0.82	0.14	44,44,45,45	0
2	ACT	A	1101	4/4	0.83	0.16	36,44,45,48	0
3	MG	A	1103	1/1	0.95	0.09	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.