



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 3ALS  
Title : Crystal structure of CEL-IV  
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Deposited on : 2010-08-07  
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

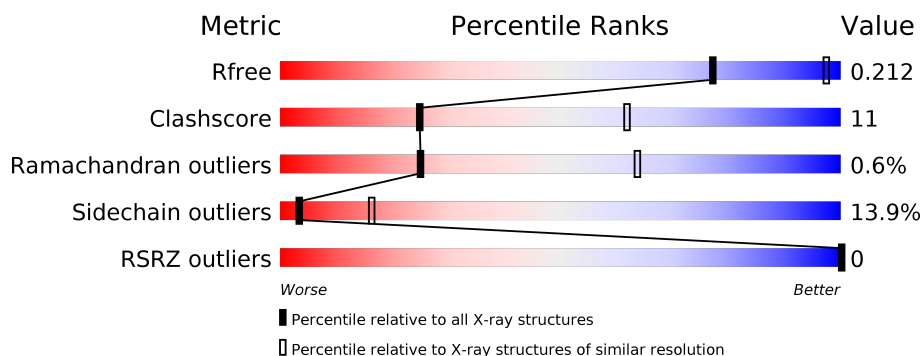
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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  $\geq 3$ , 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	157	 66% 28% 6% 0% 0%
1	B	157	 69% 26% 5% 0% 0%
1	C	157	 71% 21% 6% 2% 0%
1	D	157	 70% 23% 6% 0% 0%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4874 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 Lectin CEL-IV, C-type.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1197	752	204	234	7			
1	B	156	Total	C	N	O	S	0	0	0
			1197	752	204	234	7			
1	C	156	Total	C	N	O	S	0	0	0
			1197	752	204	234	7			
1	D	156	Total	C	N	O	S	0	0	0
			1197	752	204	234	7			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

- Molecule 3 is water.

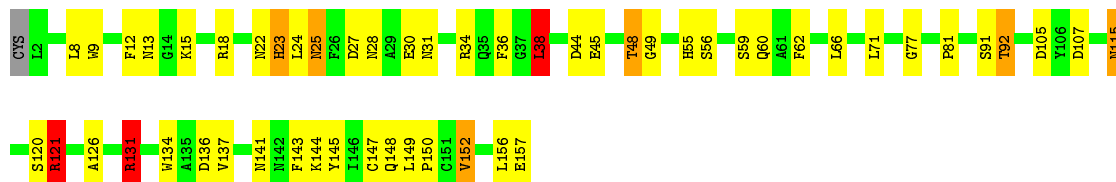
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	21	Total	O	0	0
			21	21		
3	B	18	Total	O	0	0
			18	18		
3	C	28	Total	O	0	0
			28	28		
3	D	15	Total	O	0	0
			15	15		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

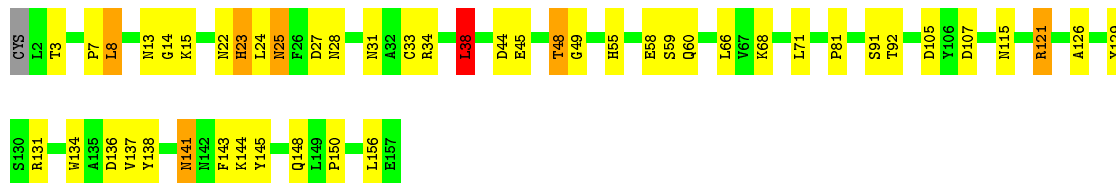
- Molecule 1: Lectin CEL-IV, C-type

Chain A: 



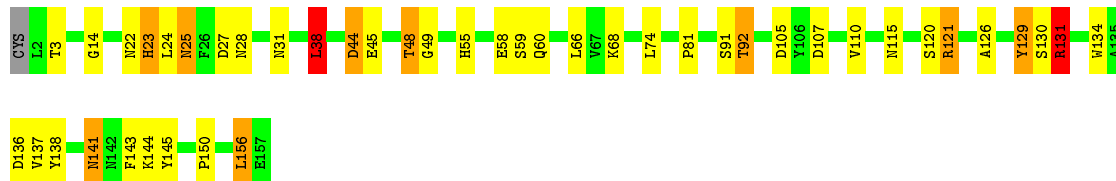
- Molecule 1: Lectin CEL-IV, C-type

Chain B: 



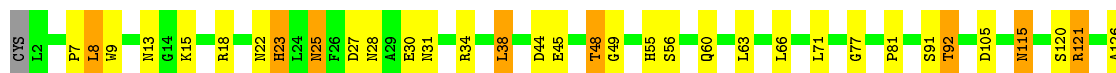
- Molecule 1: Lectin CEL-IV, C-type

Chain C: 



- Molecule 1: Lectin CEL-IV, C-type

Chain D: 



R131	W134	A135	D136	V137	N141	N142	F143	K144	Y145	L146	C147	Q148	L149	L156	E157
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.06 Å 86.06 Å 375.32 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.03 – 3.00 42.75 – 3.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.03-3.00) 100.0 (42.75-3.00)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.56 (at 3.01 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.196 , 0.215 0.195 , 0.212	Depositor DCC
$R_{free}$ test set	1575 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.3	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 8.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.489 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4874	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.97% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CA

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	1.12	3/1232 (0.2%)	1.11	10/1680 (0.6%)
1	B	1.19	5/1232 (0.4%)	1.06	5/1680 (0.3%)
1	C	1.17	5/1232 (0.4%)	1.25	8/1680 (0.5%)
1	D	1.12	2/1232 (0.2%)	1.14	8/1680 (0.5%)
All	All	1.15	15/4928 (0.3%)	1.14	31/6720 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	138	TYR	CE2-CZ	6.91	1.47	1.38
1	A	92	THR	CA-CB	6.58	1.70	1.53
1	B	138	TYR	CE2-CZ	6.56	1.47	1.38
1	D	92	THR	CA-CB	6.53	1.70	1.53
1	D	147	CYS	CB-SG	-6.49	1.71	1.82

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	131	ARG	NE-CZ-NH1	-15.09	112.75	120.30
1	C	131	ARG	NE-CZ-NH2	14.67	127.64	120.30
1	C	121	ARG	NE-CZ-NH1	-14.64	112.98	120.30
1	D	121	ARG	NE-CZ-NH1	-14.56	113.02	120.30
1	C	121	ARG	NE-CZ-NH2	13.60	127.10	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	LEU	Peptide
1	B	156	LEU	Peptide
1	C	156	LEU	Peptide
1	D	156	LEU	Peptide

## 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	1197	0	1094	29	1
1	B	1197	0	1094	22	1
1	C	1197	0	1094	25	1
1	D	1197	0	1094	26	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	21	0	0	2	0
3	B	18	0	0	0	0
3	C	28	0	0	1	0
3	D	15	0	0	0	0
All	All	4874	0	4376	99	2

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 11.

The worst 5 of 99 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:157:GLU:C	3:A:164:HOH:O	1.78	1.16
1:B:25:ASN:C	1:B:25:ASN:HD22	1.69	0.94
1:C:25:ASN:C	1:C:25:ASN:HD22	1.70	0.92
1:D:25:ASN:HD22	1:D:25:ASN:C	1.85	0.79
1:A:25:ASN:HD22	1:A:25:ASN:C	1.85	0.78

All (2) 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:77:GLY:O	1:C:129:TYR:OH[5_445]	2.08	0.12
1:B:129:TYR:OH	1:D:77:GLY:O[5_455]	2.10	0.10

## 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	154/157 (98%)	142 (92%)	11 (7%)	1 (1%)	25	64
1	B	154/157 (98%)	144 (94%)	9 (6%)	1 (1%)	25	64
1	C	154/157 (98%)	147 (96%)	6 (4%)	1 (1%)	25	64
1	D	154/157 (98%)	144 (94%)	9 (6%)	1 (1%)	25	64
All	All	616/628 (98%)	577 (94%)	35 (6%)	4 (1%)	25	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	115	ASN
1	D	115	ASN
1	B	115	ASN
1	C	115	ASN

### 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	128/129 (99%)	109 (85%)	19 (15%)	3	14
1	B	128/129 (99%)	109 (85%)	19 (15%)	3	14
1	C	128/129 (99%)	112 (88%)	16 (12%)	4	20
1	D	128/129 (99%)	111 (87%)	17 (13%)	4	17
All	All	512/516 (99%)	441 (86%)	71 (14%)	3	16

5 of 71 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	66	LEU
1	C	23	HIS
1	D	66	LEU
1	B	71	LEU
1	B	131	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	60	GLN
1	C	13	ASN
1	D	23	HIS
1	B	25	ASN
1	B	55	HIS

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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.

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	156/157 (99%)	0.01	0 100 100	25, 39, 56, 75	0
1	B	156/157 (99%)	-0.03	0 100 100	25, 39, 56, 75	0
1	C	156/157 (99%)	0.00	0 100 100	25, 39, 56, 75	0
1	D	156/157 (99%)	0.02	0 100 100	25, 39, 56, 75	0
All	All	624/628 (99%)	0.00	0 100 100	25, 39, 56, 75	0

There are no RSRZ outliers to report.

### 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](#)

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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
2	CA	D	158	1/1	0.96	0.17	32,32,32,32	0
2	CA	A	158	1/1	0.97	0.15	33,33,33,33	0
2	CA	B	158	1/1	0.99	0.16	35,35,35,35	0
2	CA	C	158	1/1	0.99	0.17	33,33,33,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.