



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 06:03 AM UTC

PDB ID : 1JWX / pdb_00001jwx
Title : Chalcone Synthase-F215S mutant
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Deposited on : 2001-09-05
Resolution : 1.63 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

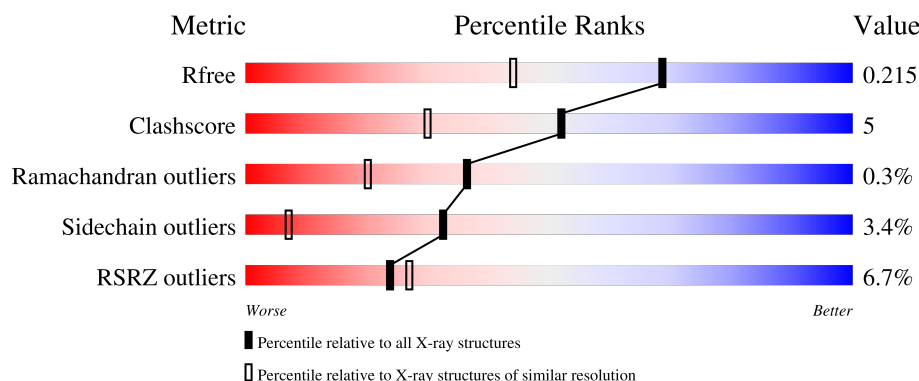
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	389	7% 85% 13% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3458 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 CHALCONE SYNTHASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	388	3114	1966	527	602	19	0	17	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	CSD	CYS	modified residue	UNP P30074
A	215	SER	PHE	engineered mutation	UNP P30074

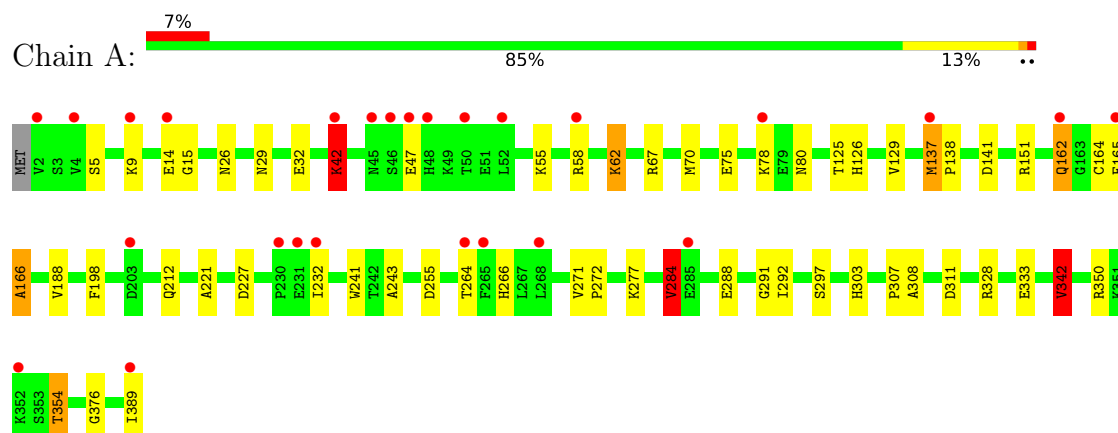
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	344	Total	O	0	0
			344	344		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: CHALCONE SYNTHASE 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.81Å 97.81Å 65.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.00 – 1.63 26.00 – 1.63	Depositor EDS
% Data completeness (in resolution range)	95.9 (26.00-1.63) 96.3 (26.00-1.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	3.50	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.29 (at 1.63Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.190 , 0.225 0.183 , 0.215	Depositor DCC
R_{free} test set	2197 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	14.9	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3458	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.28% 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: CSD

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	0.64	0/3162	1.36	18/4277 (0.4%)

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	5

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	166	ALA	N-CA-C	11.75	135.82	110.80
1	A	376	GLY	O-C-N	-11.21	108.12	122.70
1	A	232	ILE	N-CA-C	-9.56	103.34	111.56
1	A	137	MET	CA-C-O	-9.28	110.95	119.86
1	A	342	VAL	N-CA-CB	-8.72	95.73	110.65
1	A	350	ARG	NE-CZ-NH2	7.16	125.65	119.20
1	A	342	VAL	CB-CA-C	6.84	122.76	112.16
1	A	303	HIS	CA-CB-CG	-6.64	107.16	113.80
1	A	308	ALA	N-CA-CB	6.63	121.21	110.40
1	A	80	ASN	O-C-N	-6.45	115.60	121.34
1	A	284	VAL	N-CA-CB	6.12	121.84	110.77
1	A	284	VAL	CA-C-O	-6.03	113.95	120.47
1	A	138	PRO	N-CA-CB	5.96	109.16	102.60
1	A	291	GLY	CA-C-N	-5.54	115.95	123.10
1	A	291	GLY	C-N-CA	-5.54	115.95	123.10
1	A	26	ASN	O-C-N	5.46	125.21	121.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	15	GLY	CA-C-N	5.24	125.00	119.76
1	A	15	GLY	C-N-CA	5.24	125.00	119.76

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	255[B]	ASP	Mainchain
1	A	297[B]	SER	Mainchain
1	A	333[B]	GLU	Mainchain
1	A	354[B]	THR	Mainchain
1	A	42[B]	LYS	Mainchain

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	3114	0	3120	32	0
2	A	344	0	0	12	0
All	All	3458	0	3120	32	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 5.

All (32) 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:9[A]:LYS:HG3	2:A:604:HOH:O	1.89	0.72
1:A:14[A]:GLU:HG3	1:A:227:ASP:OD2	1.91	0.71
1:A:311:ASP:HB3	2:A:639:HOH:O	1.92	0.70
1:A:307:PRO:HD2	2:A:733:HOH:O	1.99	0.62
1:A:264:THR:HG22	2:A:570:HOH:O	2.01	0.60
1:A:58:ARG:HG2	1:A:62:LYS:HD2	1.83	0.60
1:A:307:PRO:CD	2:A:733:HOH:O	2.51	0.59
1:A:264:THR:HG23	1:A:266:HIS:NE2	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ARG:NH2	1:A:62:LYS:HE3	2.20	0.56
1:A:165:PHE:CD1	1:A:165:PHE:C	2.84	0.55
1:A:126:HIS:HE1	2:A:425:HOH:O	1.90	0.54
1:A:58:ARG:HH21	1:A:62:LYS:HE3	1.73	0.53
1:A:125:THR:OG1	1:A:126:HIS:HD2	1.93	0.50
1:A:151:ARG:HA	1:A:151:ARG:HE	1.76	0.49
1:A:55:LYS:HE3	2:A:725:HOH:O	2.13	0.48
1:A:14[A]:GLU:CG	1:A:227:ASP:OD2	2.59	0.48
1:A:42[B]:LYS:HG3	1:A:47:GLU:CD	2.40	0.46
1:A:29:ASN:HB3	1:A:70:MET:O	2.15	0.46
1:A:188:VAL:O	1:A:221:ALA:HA	2.16	0.46
1:A:198:PHE:HA	1:A:212:GLN:HE22	1.83	0.44
1:A:277:LYS:HG3	2:A:681:HOH:O	2.18	0.43
1:A:162:GLN:HB2	2:A:515:HOH:O	2.18	0.43
1:A:137:MET:HE3	1:A:137:MET:HB2	1.79	0.42
1:A:271:VAL:HB	1:A:272:PRO:HD3	2.00	0.42
1:A:32[2]:GLU:HG3	2:A:656:HOH:O	2.20	0.41
1:A:129:VAL:HG21	1:A:141:ASP:HA	2.03	0.41
1:A:342:VAL:HB	2:A:399:HOH:O	2.21	0.40
1:A:241:TRP:CH2	1:A:243:ALA:HB2	2.56	0.40
1:A:284:VAL:O	1:A:288:GLU:HB2	2.21	0.40
1:A:328[A]:ARG:HD2	2:A:649:HOH:O	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	401/389 (103%)	391 (98%)	9 (2%)	1 (0%)	43 27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	ALA

5.3.2 Protein sidechains [i](#)

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	339/324 (105%)	326 (96%)	13 (4%)	29 5

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5[A]	SER
1	A	5[B]	SER
1	A	42[A]	LYS
1	A	42[B]	LYS
1	A	62	LYS
1	A	67	ARG
1	A	75	GLU
1	A	78	LYS
1	A	162	GLN
1	A	284	VAL
1	A	292	ILE
1	A	342	VAL
1	A	389	ILE

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

Mol	Chain	Res	Type
1	A	126	HIS
1	A	212	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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	$\# Z > 2$
1	CSD	A	164	1	4,7,8	8.66	1 (25%)	1,8,10	4.95	1 (100%)

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
1	CSD	A	164	1	-	0/2/6/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	164	CSD	OD1-SG	-17.23	1.32	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	CSD	OD1-SG-CB	-4.95	96.49	105.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no oligosaccharides 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	387/389 (99%)	0.29	26 (6%)	24 27	8, 16, 29, 42	16 (4%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	389	ILE	4.9
1	A	50	THR	3.9
1	A	264	THR	3.8
1	A	2	VAL	3.2
1	A	48	HIS	3.1
1	A	9[A]	LYS	3.1
1	A	231	GLU	3.1
1	A	268	LEU	2.9
1	A	352	LYS	2.8
1	A	232	ILE	2.7
1	A	47	GLU	2.6
1	A	52	LEU	2.5
1	A	165	PHE	2.5
1	A	58	ARG	2.4
1	A	78	LYS	2.3
1	A	45	ASN	2.2
1	A	285	GLU	2.2
1	A	203	ASP	2.2
1	A	46	SER	2.2
1	A	162	GLN	2.1
1	A	42[A]	LYS	2.1
1	A	4	VAL	2.1
1	A	230	PRO	2.1
1	A	14[A]	GLU	2.0
1	A	137	MET	2.0
1	A	265	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSD	A	164	8/9	0.75	0.23	17,23,28,30	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.