



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 09:58 AM GMT

PDB ID : 3KF5
Title : Structure of invertase from Schwanniomyces occidentalis
Authors : Sanz-Aparicio, J.; Polo, A.
Deposited on : 2009-10-27
Resolution : 2.90 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

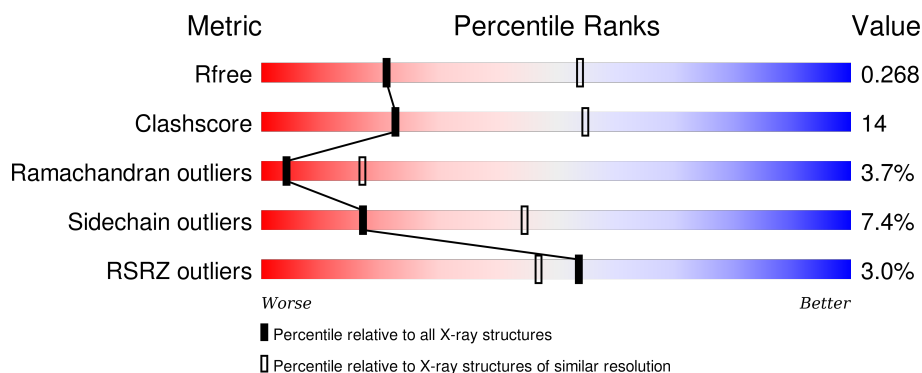
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	512	<div> <div>3%</div> <div>70%</div> <div>25%</div> <div>• •</div> </div>
1	B	512	<div> <div>3%</div> <div>70%</div> <div>24%</div> <div>• ••</div> </div>

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	NAG	A	5001	-	-	-	X

2 Entry composition [i](#)

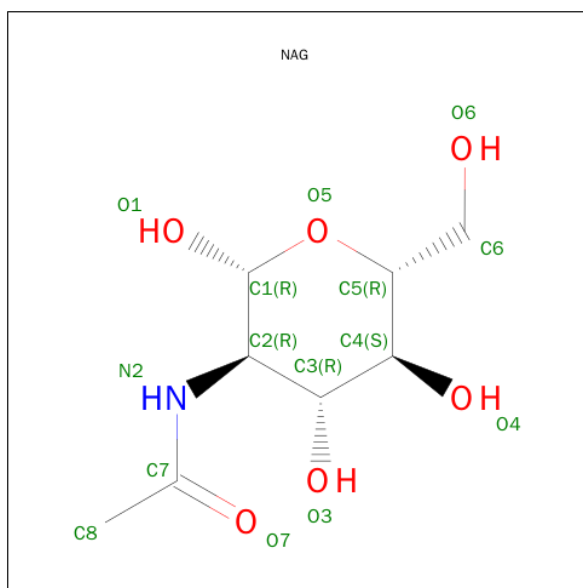
There are 4 unique types of molecules in this entry. The entry contains 8626 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Invertase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	512	Total	C	N	O	S	0	0	0
			4156	2670	680	800	6			
1	B	508	Total	C	N	O	S	0	0	0
			4126	2651	676	793	6			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



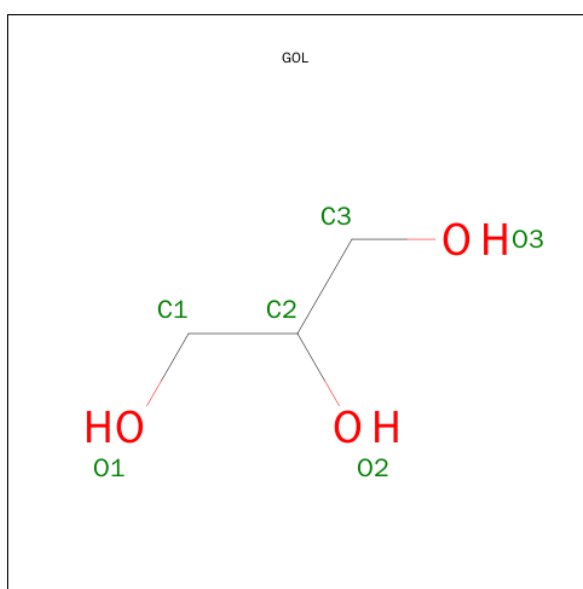
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	119	Total	O	0	0
			119	119		
4	B	95	Total	O	0	0
			95	95		

1485	1486	1487	1488	1495	1496	1497	F502	M503	H510	B511	I512	Q513	T516	D517	T518	F523	I529	L533	M534	K535
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	105.78Å 119.49Å 137.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.65 – 2.90 59.65 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (59.65-2.90) 99.8 (59.65-2.90)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.222 , 0.257 0.228 , 0.268	Depositor DCC
R_{free} test set	1979 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 28.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	2 of 39315 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	8626	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, NAG

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	0.43	0/4277	0.66	2/5825 (0.0%)
1	B	0.42	0/4247	0.65	2/5784 (0.0%)
All	All	0.43	0/8524	0.66	4/11609 (0.0%)

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	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	463	LEU	N-CA-C	5.76	126.55	111.00
1	A	463	LEU	N-CA-C	5.53	125.92	111.00
1	A	502	PHE	N-CA-C	5.31	125.33	111.00
1	B	502	PHE	N-CA-C	5.04	124.59	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	359	ASP	Peptide
1	A	459	LEU	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	B	359	ASP	Peptide
1	B	459	LEU	Peptide

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	4156	0	3937	117	0
1	B	4126	0	3908	110	0
2	A	70	0	65	1	0
2	B	42	0	39	1	0
3	A	18	0	24	1	0
4	A	119	0	0	2	0
4	B	95	0	0	6	0
All	All	8626	0	7973	222	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 14.

The worst 5 of 222 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:B:116:VAL:HG11	1:B:132:ASN:HA	1.42	0.99
1:A:116:VAL:HG11	1:A:132:ASN:HA	1.43	0.97
1:B:36:ARG:HD2	1:B:42:THR:HG21	1.47	0.94
1:B:170:ILE:HD12	1:B:193:MET:HE1	1.49	0.93
1:A:371:ASN:HD22	1:A:371:ASN:N	1.62	0.93

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	510/512 (100%)	446 (88%)	44 (9%)	20 (4%)	4	15
1	B	506/512 (99%)	447 (88%)	41 (8%)	18 (4%)	4	18
All	All	1016/1024 (99%)	893 (88%)	85 (8%)	38 (4%)	4	17

5 of 38 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	26	ASP
1	A	127	SER
1	A	187	ASP
1	A	303	HIS
1	A	360	SER

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	464/464 (100%)	428 (92%)	36 (8%)	16	41
1	B	460/464 (99%)	428 (93%)	32 (7%)	19	47
All	All	924/928 (100%)	856 (93%)	68 (7%)	17	44

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	470	GLN
1	B	44	GLU
1	B	470	GLN
1	A	483	LYS
1	A	495	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	470	GLN
1	B	52	ASN
1	B	453	GLN
1	A	484	ASN
1	A	510	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 ligands are 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	$\# Z > 2$
2	NAG	A	1001	1	14,14,15	0.52	0	15,19,21	1.78	2 (13%)
2	NAG	A	2001	1	14,14,15	0.42	0	15,19,21	1.16	1 (6%)
2	NAG	A	3001	1	14,14,15	0.63	0	15,19,21	1.78	1 (6%)
2	NAG	A	4001	1	14,14,15	0.43	0	15,19,21	1.58	2 (13%)
2	NAG	A	5001	1	14,14,15	0.64	0	15,19,21	1.69	3 (20%)
3	GOL	A	6001	-	5,5,5	0.38	0	5,5,5	0.13	0
3	GOL	A	6002	-	5,5,5	0.29	0	5,5,5	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	6003	-	5,5,5	0.34	0	5,5,5	0.38	0
2	NAG	B	3001	1	14,14,15	0.59	0	15,19,21	2.36	2 (13%)
2	NAG	B	4001	1	14,14,15	0.59	0	15,19,21	1.43	3 (20%)
2	NAG	B	5001	1	14,14,15	0.59	0	15,19,21	2.20	3 (20%)

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	NAG	A	1001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	2001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	3001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	4001	1	-	0/6/23/26	0/1/1/1
2	NAG	A	5001	1	-	0/6/23/26	0/1/1/1
3	GOL	A	6001	-	-	0/4/4/4	0/0/0/0
3	GOL	A	6002	-	-	0/4/4/4	0/0/0/0
3	GOL	A	6003	-	-	0/4/4/4	0/0/0/0
2	NAG	B	3001	1	-	0/6/23/26	0/1/1/1
2	NAG	B	4001	1	-	0/6/23/26	0/1/1/1
2	NAG	B	5001	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4001	NAG	C3-C4-C5	-2.96	105.04	110.20
2	B	3001	NAG	C6-C5-C4	-2.62	106.55	113.02
2	B	5001	NAG	C3-C2-N2	-2.58	104.39	110.56
2	A	5001	NAG	C3-C2-N2	-2.39	104.83	110.56
2	B	4001	NAG	O4-C4-C5	2.37	115.53	109.24

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	4001	NAG	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	6001	GOL	1	0
2	B	3001	NAG	1	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	512/512 (100%)	0.17	15 (2%) 55 49	10, 19, 33, 46	0
1	B	508/512 (99%)	0.22	16 (3%) 52 45	10, 19, 32, 46	0
All	All	1020/1024 (99%)	0.20	31 (3%) 54 47	10, 19, 32, 46	0

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	128	SER	4.6
1	B	409	THR	4.2
1	A	127	SER	3.9
1	A	24	SER	3.8
1	A	26	ASP	3.4

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	A	5001	14/15	0.91	0.29	8.17	37,43,46,46	0
3	GOL	A	6001	6/6	0.95	0.19	1.08	31,35,37,38	0
2	NAG	A	3001	14/15	0.89	0.21	0.99	30,34,38,39	0
2	NAG	B	3001	14/15	0.87	0.21	0.77	31,36,38,38	0
3	GOL	A	6002	6/6	0.91	0.19	0.54	59,60,61,61	0
2	NAG	B	5001	14/15	0.87	0.21	0.32	37,42,44,44	0
2	NAG	B	4001	14/15	0.83	0.26	0.29	30,36,37,37	0
2	NAG	A	1001	14/15	0.91	0.21	-0.03	36,39,42,43	0
2	NAG	A	4001	14/15	0.91	0.18	-0.68	30,34,36,37	0
2	NAG	A	2001	14/15	0.84	0.47	-	51,57,59,59	0
3	GOL	A	6003	6/6	0.80	0.29	-	51,52,52,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.