



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:41 PM GMT

PDB ID : 1Q4L  
Title : GSK-3 Beta complexed with Inhibitor I-5  
Authors : Bertrand, J.A.; Thieffine, S.; Vulpetti, A.; Cristiani, C.; Valsasina, B.; Knapp, S.; Kalisz, H.M.; Flocco, M.  
Deposited on : 2003-08-04  
Resolution : 2.77 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  
<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

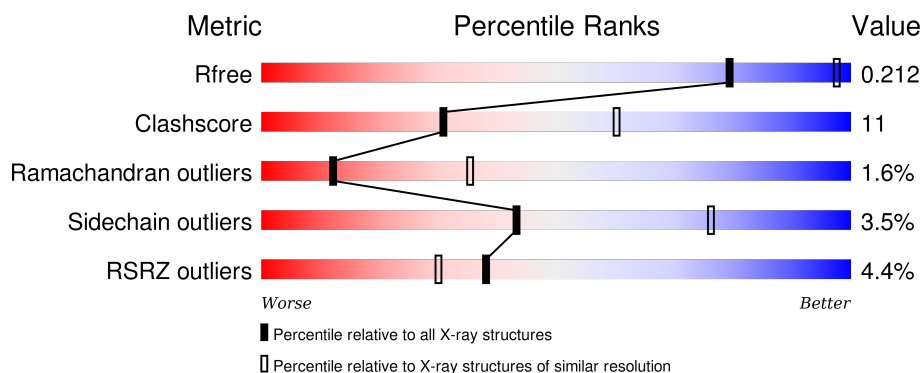
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.77 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	3004 (2.80-2.76)
Clashscore	102246	3480 (2.80-2.76)
Ramachandran outliers	100387	3423 (2.80-2.76)
Sidechain outliers	100360	3425 (2.80-2.76)
RSRZ outliers	91569	3016 (2.80-2.76)

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. 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	424	<div> <div>4%</div> <div>61%</div> <div>18%</div> <div>•</div> <div>19%</div> </div>
1	B	424	<div> <div>3%</div> <div>60%</div> <div>17%</div> <div>•</div> <div>21%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5515 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.

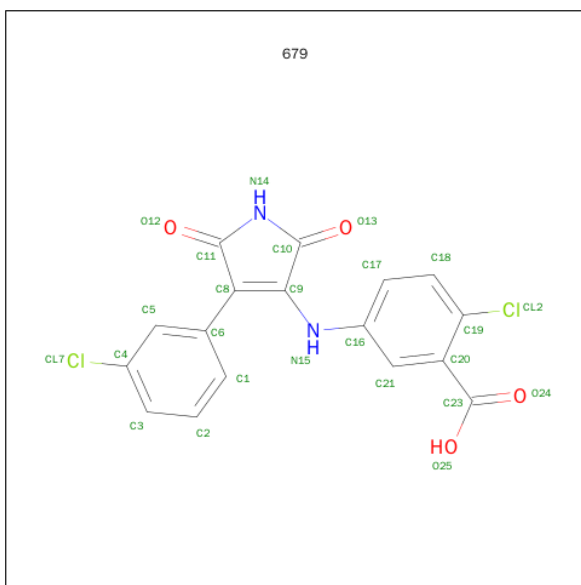
- Molecule 1 is a protein called GLYCOGEN SYNTHASE KINASE-3 BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2728	1753	470	494	11			
1	B	337	Total	C	N	O	S	0	0	0
			2690	1726	461	492	11			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	CLONING ARTIFACT	UNP P49841
A	-2	PRO	-	CLONING ARTIFACT	UNP P49841
A	-1	LEU	-	CLONING ARTIFACT	UNP P49841
A	0	GLY	-	CLONING ARTIFACT	UNP P49841
A	1	SER	-	CLONING ARTIFACT	UNP P49841
B	-3	GLY	-	CLONING ARTIFACT	UNP P49841
B	-2	PRO	-	CLONING ARTIFACT	UNP P49841
B	-1	LEU	-	CLONING ARTIFACT	UNP P49841
B	0	GLY	-	CLONING ARTIFACT	UNP P49841
B	1	SER	-	CLONING ARTIFACT	UNP P49841

- Molecule 2 is 2-CHLORO-5-[4-(3-CHLORO-PHENYL)-2,5-DIOXO-2,5-DIHYDRO-1H-PYRROL-3-YLAMINO]-BENZOIC ACID (three-letter code: 679) (formula: C<sub>17</sub>H<sub>10</sub>Cl<sub>2</sub>N<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			25	17	2	2	4		
2	B	1	Total	C	Cl	N	O	0	0
			25	17	2	2	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		
3	B	30	Total	O	0	0
			30	30		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.17Å 86.66Å 178.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.77 20.00 – 2.77	Depositor EDS
% Data completeness (in resolution range)	98.7 (20.00-2.77) 98.8 (20.00-2.77)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.79Å)	Xtriage
Refinement program	CNX 2000	Depositor
R, $R_{free}$	0.212 , 0.251 0.208 , 0.212	Depositor DCC
$R_{free}$ test set	1704 reflections (5.36%)	DCC
Wilson B-factor (Å <sup>2</sup> )	50.5	Xtriage
Anisotropy	0.596	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 52.2	EDS
Estimated twinning fraction	0.023 for k,h,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 33498 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5515	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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, 0.375 respectively for untwinned datasets, and 0.333, 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:  
679

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	0.45	0/2794	0.67	0/3800
1	B	0.47	0/2755	0.68	0/3748
All	All	0.46	0/5549	0.68	0/7548

There are no bond length outliers.

There are no bond angle outliers.

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	2728	0	2755	61	0
1	B	2690	0	2707	58	0
2	A	25	0	9	1	0
2	B	25	0	9	2	0
3	A	17	0	0	0	0
3	B	30	0	0	0	0
All	All	5515	0	5480	119	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 11.

The worst 5 of 119 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:383:ARG:HE	1:B:383:ARG:HA	1.08	1.08
1:B:89:GLN:HE22	1:B:95:ASN:HB2	1.38	0.89
1:B:383:ARG:NE	1:B:383:ARG:HA	1.89	0.87
1:B:383:ARG:HE	1:B:383:ARG:CA	1.89	0.85
1:B:62:ILE:HG21	1:B:72:GLN:HB2	1.64	0.79

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	336/424 (79%)	311 (93%)	17 (5%)	8 (2%)	7	23
1	B	331/424 (78%)	309 (93%)	19 (6%)	3 (1%)	21	53
All	All	667/848 (79%)	620 (93%)	36 (5%)	11 (2%)	12	35

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	ALA
1	B	66	SER
1	B	125	GLU
1	A	49	ASP
1	A	221	TYR

### 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	303/366 (83%)	294 (97%)	9 (3%)	48	81
1	B	300/366 (82%)	288 (96%)	12 (4%)	38	72
All	All	603/732 (82%)	582 (96%)	21 (4%)	43	76

5 of 21 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	90	ASP
1	B	101	MET
1	B	316	LEU
1	A	370	ASN
1	B	370	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	370	ASN
1	B	370	ASN
1	B	89	GLN
1	A	151	GLN
1	B	72	GLN

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

2 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	679	A	451	-	23,27,27	2.11	12 (52%)	30,39,39	1.77	6 (20%)
2	679	B	452	-	23,27,27	2.22	13 (56%)	30,39,39	1.86	8 (26%)

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	679	A	451	-	-	0/8/28/28	0/3/3/3
2	679	B	452	-	-	0/8/28/28	0/3/3/3

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	452	679	C8-C11	-3.47	1.42	1.50
2	A	451	679	C6-C8	-3.36	1.43	1.49
2	A	451	679	C8-C11	-3.09	1.42	1.50
2	B	452	679	C6-C8	-2.25	1.45	1.49
2	B	452	679	C10-N14	-2.15	1.33	1.37

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	679	C18-C19-CL2	-3.06	112.06	118.39
2	B	452	679	C9-C8-C11	-2.29	103.50	106.51
2	B	452	679	C18-C19-CL2	-2.20	113.83	118.39
2	A	451	679	C11-N14-C10	-2.12	109.17	111.29
2	B	452	679	C20-C19-CL2	2.05	123.82	120.67

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	451	679	1	0
2	B	452	679	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, 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	342/424 (80%)	-0.06	18 (5%)	30 22	25, 48, 109, 170	0
1	B	337/424 (79%)	-0.22	12 (3%)	46 39	23, 44, 101, 157	0
All	All	679/848 (80%)	-0.14	30 (4%)	38 30	23, 46, 108, 170	0

The worst 5 of 30 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	93	PHE	7.7
1	A	90	ASP	4.9
1	A	295	GLN	4.7
1	B	278	ARG	4.2
1	A	385	GLN	3.9

### 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, 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	679	A	451	25/25	0.92	0.18	0.78	47,47,47,47	0
2	679	B	452	25/25	0.93	0.16	-0.06	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.