



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 18, 2024 – 03:22 AM EDT

PDB ID : 5VE6
Title : Crystal structure of Sugren kinase 223
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Deposited on : 2017-04-03
Resolution : 2.95 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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
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)	:	2.37.1

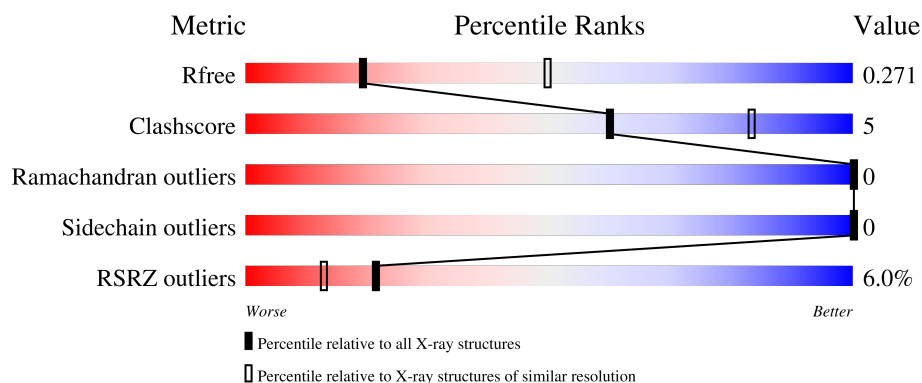
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.95 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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	492	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2655 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 Tyrosine-protein kinase SgK223.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	348	Total	C	N	O	S	0	0	0
			2646	1686	452	485	23			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	913	MET	-	initiating methionine	UNP Q86YV5
A	914	HIS	-	expression tag	UNP Q86YV5
A	915	HIS	-	expression tag	UNP Q86YV5
A	916	HIS	-	expression tag	UNP Q86YV5
A	917	HIS	-	expression tag	UNP Q86YV5
A	918	HIS	-	expression tag	UNP Q86YV5
A	919	HIS	-	expression tag	UNP Q86YV5
A	920	HIS	-	expression tag	UNP Q86YV5
A	921	HIS	-	expression tag	UNP Q86YV5
A	922	GLY	-	expression tag	UNP Q86YV5
A	923	SER	-	expression tag	UNP Q86YV5
A	924	GLU	-	expression tag	UNP Q86YV5
A	925	ASN	-	expression tag	UNP Q86YV5
A	926	LEU	-	expression tag	UNP Q86YV5
A	927	TYR	-	expression tag	UNP Q86YV5
A	928	PHE	-	expression tag	UNP Q86YV5
A	929	GLN	-	expression tag	UNP Q86YV5
A	930	GLY	-	expression tag	UNP Q86YV5
A	931	SER	-	expression tag	UNP Q86YV5
A	?	-	PRO	deletion	UNP Q86YV5
A	?	-	ALA	deletion	UNP Q86YV5

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		

- Molecule 1: Tyrosine-protein kinase SgK223



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.68Å 144.68Å 47.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.75 – 2.95 45.75 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.75-2.95) 99.8 (45.75-2.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.37 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.218 , 0.262 0.229 , 0.271	Depositor DCC
R_{free} test set	955 reflections (8.63%)	wwPDB-VP
Wilson B-factor (Å ²)	65.7	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2655	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.23	0/2698	0.39	0/3665

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 [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	2646	0	2556	24	1
2	A	9	0	0	0	0
All	All	2655	0	2556	24	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 5.

All (24) 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:1016:ASP:OD1	1:A:1019:SER:OG	2.12	0.64
1:A:1335:GLU:N	1:A:1335:GLU:OE1	2.31	0.63
1:A:1051:CYS:SG	1:A:1093:ARG:NH1	2.74	0.60
1:A:1209:LEU:HD12	1:A:1210:PRO:HD2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:948:LYS:O	1:A:951:THR:OG1	2.17	0.59
1:A:1327:CYS:HA	1:A:1331:GLY:HA3	1.87	0.55
1:A:1092:THR:HG22	1:A:1094:GLU:H	1.74	0.53
1:A:1042:PRO:HG2	1:A:1139:HIS:CD2	2.45	0.52
1:A:1003:SER:HB2	1:A:1149:GLU:O	2.14	0.48
1:A:1283:ARG:HG3	1:A:1283:ARG:HH11	1.78	0.48
1:A:1317:ARG:HG3	1:A:1320:GLU:OE2	2.13	0.48
1:A:1309:GLU:O	1:A:1315:ARG:NH1	2.47	0.47
1:A:1251:LYS:HB3	1:A:1315:ARG:HE	1.80	0.47
1:A:1133:LEU:HD13	1:A:1253:ASP:HB3	1.98	0.46
1:A:1316:ILE:HG23	1:A:1320:GLU:HB2	1.97	0.46
1:A:947:SER:O	1:A:951:THR:HG23	2.17	0.45
1:A:1144:ARG:O	1:A:1252:PHE:HE2	2.00	0.44
1:A:1243:ILE:HG23	1:A:1244:VAL:HG23	2.00	0.44
1:A:1311:ASP:HA	1:A:1312:PRO:HD3	1.85	0.43
1:A:961:GLN:HG3	1:A:965:ARG:HH22	1.83	0.42
1:A:1265:LEU:HD11	1:A:1300:LEU:HD13	2.01	0.41
1:A:1251:LYS:HG3	1:A:1312:PRO:O	2.20	0.41
1:A:1054:PHE:HE1	1:A:1056:ALA:HB2	1.85	0.41
1:A:1297:SER:HB2	1:A:1298:PRO:HD3	2.03	0.41

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)
1:A:959:TYR:OH	1:A:959:TYR:OH[7_555]	2.02	0.18

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	334/492 (68%)	325 (97%)	9 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	278/415 (67%)	278 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	961	GLN
1	A	1150	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

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 [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	348/492 (70%)	0.62	21 (6%) 21 13	35, 57, 83, 100	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	951	THR	4.8
1	A	1019	SER	3.5
1	A	1082	ALA	3.4
1	A	1348	CYS	3.4
1	A	1158	LEU	3.0
1	A	947	SER	3.0
1	A	1393	PRO	2.8
1	A	1016	ASP	2.7
1	A	950	GLY	2.7
1	A	955	LEU	2.6
1	A	989	TRP	2.6
1	A	956	GLY	2.5
1	A	1002	ASP	2.4
1	A	959	TYR	2.4
1	A	1087	CYS	2.3
1	A	958	LEU	2.3
1	A	1405	LEU	2.2
1	A	1091	ILE	2.1
1	A	1391	ALA	2.1
1	A	1249	TYR	2.1
1	A	1366	PHE	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 monosaccharides 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.