WORLDWIDE PDB

PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report (i)

Sep 3, 2023 – 01:04 PM EDT

PDB ID	:	3SJB
Title	:	Crystal structure of S. cerevisiae Get3 in the open state in complex with Get1 cytosolic domain
Authors	:	Reitz, S.; Wild, K.; Sinning, I.
Deposited on	:	2011-06-21
Resolution	:	3.30 Å(reported)

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

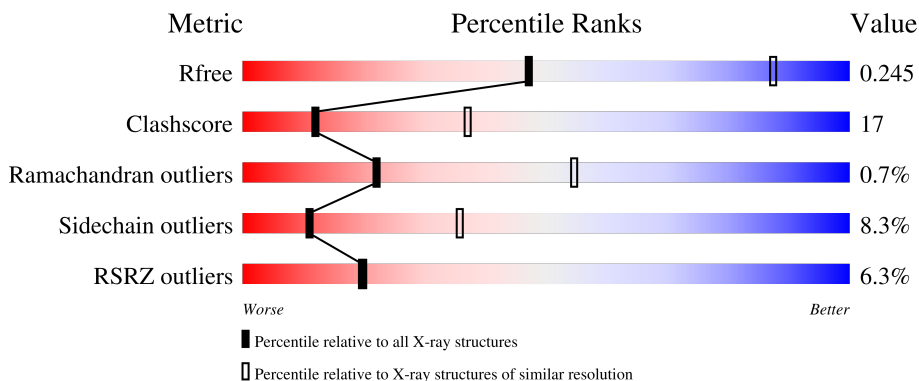|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 3.30 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	Percentile Ranks	Value
Rfree		0.245
Clashscore		17
Ramachandran outliers		0.7%
Sidechain outliers		8.3%
RSRZ outliers		6.3%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 <=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	362	4% 50% / 26% / • / 20%
1	B	362	5% 52% / 28% / • / 17%
2	C	93	6% 49% / 17% / • / 32%
2	D	93	10% 48% / 27% / • / 23%

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 5823 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 ATPase GET3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2280	1437	380	447	16			
1	B	302	Total	C	N	O	S	0	0	0
			2384	1503	396	468	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	355	LEU	-	expression tag	UNP Q12154
A	356	GLU	-	expression tag	UNP Q12154
A	357	HIS	-	expression tag	UNP Q12154
A	358	HIS	-	expression tag	UNP Q12154
A	359	HIS	-	expression tag	UNP Q12154
A	360	HIS	-	expression tag	UNP Q12154
A	361	HIS	-	expression tag	UNP Q12154
A	362	HIS	-	expression tag	UNP Q12154
B	355	LEU	-	expression tag	UNP Q12154
B	356	GLU	-	expression tag	UNP Q12154
B	357	HIS	-	expression tag	UNP Q12154
B	358	HIS	-	expression tag	UNP Q12154
B	359	HIS	-	expression tag	UNP Q12154
B	360	HIS	-	expression tag	UNP Q12154
B	361	HIS	-	expression tag	UNP Q12154
B	362	HIS	-	expression tag	UNP Q12154

- Molecule 2 is a protein called Golgi to ER traffic protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	63	Total	C	N	O	0	0	0
			528	328	96	104			
2	D	72	Total	C	N	O	0	0	0
			610	379	115	116			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	18	MET	-	expression tag	UNP P53192
C	104	GLU	-	expression tag	UNP P53192
C	105	HIS	-	expression tag	UNP P53192
C	106	HIS	-	expression tag	UNP P53192
C	107	HIS	-	expression tag	UNP P53192
C	108	HIS	-	expression tag	UNP P53192
C	109	HIS	-	expression tag	UNP P53192
C	110	HIS	-	expression tag	UNP P53192
D	18	MET	-	expression tag	UNP P53192
D	104	GLU	-	expression tag	UNP P53192
D	105	HIS	-	expression tag	UNP P53192
D	106	HIS	-	expression tag	UNP P53192
D	107	HIS	-	expression tag	UNP P53192
D	108	HIS	-	expression tag	UNP P53192
D	109	HIS	-	expression tag	UNP P53192
D	110	HIS	-	expression tag	UNP P53192

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).
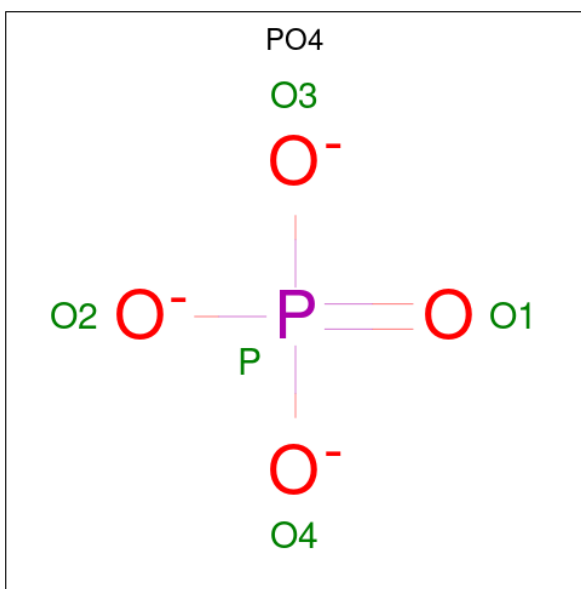
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 5	O 4	P 1	0	0
3	A	1	Total 5	O 4	P 1	0	0
3	B	1	Total 5	O 4	P 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total 5	O 4	P 1	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Zn 1	0	0

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: ATPase GET3

• Molecule 1: ATPase GET3

• Molecule 2: Golgi to ER traffic protein 1

Chain C:

• Molecule 2: Golgi to ER traffic protein 1

Chain D:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α, β, γ	112.90Å 112.90Å 310.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.89 – 3.30 48.89 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.89-3.30) 99.6 (48.89-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.27 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6_289)	Depositor
R, R_{free}	0.195 , 0.243 0.193 , 0.245	Depositor DCC
R_{free} test set	941 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	110.7	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 103.5	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5823	wwPDB-VP
Average B, all atoms (Å^2)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: PO4, ZN

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	0.47	0/2318	0.63	0/3125
1	B	0.45	0/2423	0.61	0/3268
2	C	0.39	0/535	0.47	0/713
2	D	0.35	0/621	0.48	0/829
All	All	0.45	0/5897	0.60	0/7935

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	2280	0	2246	82	0
1	B	2384	0	2357	91	0
2	C	528	0	523	10	0
2	D	610	0	598	19	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
4	A	1	0	0	0	0
All	All	5823	0	5724	198	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 17.

The worst 5 of 198 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:20:TRP:HB2	1:B:236:THR:HB	1.45	0.98
1:A:20:TRP:HB2	1:A:236:THR:HB	1.47	0.96
1:A:29:VAL:HG12	1:A:243:ILE:HD13	1.44	0.96
1:B:29:VAL:HG12	1:B:243:ILE:HD13	1.44	0.95
1:A:72:LYS:N	1:A:72:LYS:HD3	1.87	0.87

There are no symmetry-related clashes.

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	282/362 (78%)	254 (90%)	26 (9%)	2 (1%)	22	54
1	B	296/362 (82%)	272 (92%)	22 (7%)	2 (1%)	22	54
2	C	61/93 (66%)	57 (93%)	4 (7%)	0	100	100
2	D	70/93 (75%)	66 (94%)	3 (4%)	1 (1%)	11	38
All	All	709/910 (78%)	649 (92%)	55 (8%)	5 (1%)	22	54

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	283	HIS
1	B	60	HIS
1	B	283	HIS
2	D	103	LEU

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	257/317 (81%)	233 (91%)	24 (9%)	9	30
1	B	269/317 (85%)	243 (90%)	26 (10%)	8	29
2	C	58/86 (67%)	56 (97%)	2 (3%)	37	65
2	D	67/86 (78%)	65 (97%)	2 (3%)	41	68
All	All	651/806 (81%)	597 (92%)	54 (8%)	11	36

5 of 54 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	80	MET
1	B	236	THR
1	B	354	GLU
1	B	81	ASN
1	B	144	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	36	ASN
1	B	326	ASN
1	B	61	ASN
1	B	289	GLN
1	B	18	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 ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	$\#|Z| > 2$
3	PO4	A	363	-	4,4,4	0.80	0	6,6,6	0.76	0
3	PO4	B	364	-	4,4,4	0.81	0	6,6,6	0.64	0
3	PO4	B	363	-	4,4,4	0.84	0	6,6,6	0.41	0
3	PO4	A	364	-	4,4,4	0.87	0	6,6,6	0.58	0

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 ⓘ

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	288/362 (79%)	0.14	13 (4%)	33	32	65, 109, 218, 258	0
1	B	302/362 (83%)	0.23	18 (5%)	21	21	64, 111, 223, 268	0
2	C	63/93 (67%)	0.35	6 (9%)	8	8	86, 129, 202, 219	0
2	D	72/93 (77%)	0.40	9 (12%)	3	3	87, 139, 229, 256	0
All	All	725/910 (79%)	0.22	46 (6%)	20	20	64, 114, 221, 268	0

The worst 5 of 46 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	99	ASP	6.1
1	B	128	ASP	6.1
1	B	223	VAL	5.7
1	B	219	LEU	5.0
1	B	184	SER	4.7

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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(Å^2)	Q<0.9
3	PO4	B	364	5/5	0.84	0.28	122,143,153,179	0
3	PO4	A	364	5/5	0.85	0.41	140,151,158,175	0
3	PO4	B	363	5/5	0.93	0.37	166,169,178,184	0
3	PO4	A	363	5/5	0.93	0.35	150,152,160,172	0
4	ZN	A	365	1/1	1.00	0.10	106,106,106,106	0

6.5 Other polymers ⓘ

There are no such residues in this entry.