



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:36 PM UTC

PDB ID : 3KXK / pdb_00003kxk
Title : Crystal structure of SsGBP mutation variant G235P
Authors : Huang, B.; Li, X.; Zhang, X.C.; Rao, Z.
Deposited on : 2009-12-03
Resolution : 2.35 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
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

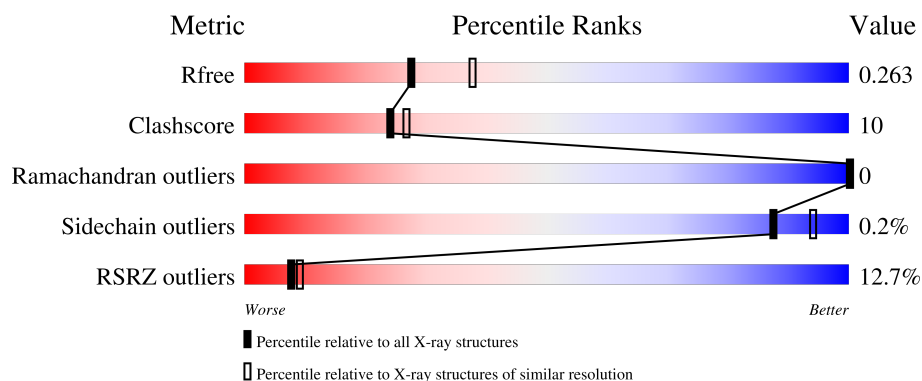
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	364	7% 63% 18% 20%
1	B	364	13% 57% 21% 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4720 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 GTP-binding protein (HflX).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	35	0	0
			2360	1538	385	434	3			
1	B	286	Total	C	N	O	S	98	0	0
			2295	1499	371	423	2			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	235	PRO	GLY	engineered mutation	UNP Q980M3
A	357	LEU	-	expression tag	UNP Q980M3
A	358	GLU	-	expression tag	UNP Q980M3
A	359	HIS	-	expression tag	UNP Q980M3
A	360	HIS	-	expression tag	UNP Q980M3
A	361	HIS	-	expression tag	UNP Q980M3
A	362	HIS	-	expression tag	UNP Q980M3
A	363	HIS	-	expression tag	UNP Q980M3
A	364	HIS	-	expression tag	UNP Q980M3
B	235	PRO	GLY	engineered mutation	UNP Q980M3
B	357	LEU	-	expression tag	UNP Q980M3
B	358	GLU	-	expression tag	UNP Q980M3
B	359	HIS	-	expression tag	UNP Q980M3
B	360	HIS	-	expression tag	UNP Q980M3
B	361	HIS	-	expression tag	UNP Q980M3
B	362	HIS	-	expression tag	UNP Q980M3
B	363	HIS	-	expression tag	UNP Q980M3
B	364	HIS	-	expression tag	UNP Q980M3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	21	Total	O	0	0
			21	21		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.71Å 71.24Å 97.67Å 90.00° 96.81° 90.00°	Depositor
Resolution (Å)	38.82 – 2.35 38.82 – 2.35	Depositor EDS
% Data completeness (in resolution range)	95.0 (38.82-2.35) 95.0 (38.82-2.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.244 , 0.266 0.241 , 0.263	Depositor DCC
R_{free} test set	1567 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4720	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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

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	0.26	0/2397	0.67	0/3231
1	B	0.26	0/2331	0.68	0/3145
All	All	0.26	0/4728	0.67	0/6376

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	2360	0	2490	48	0
1	B	2295	0	2434	48	0
2	A	44	0	0	0	0
2	B	21	0	0	1	0
All	All	4720	0	4924	93	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 10.

All (93) 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:113:LYS:O	1:A:117:PRO:HD2	1.75	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ILE:HD12	1:A:66:LEU:HD11	1.66	0.78
1:B:9:SER:HB3	1:B:33:LEU:HD22	1.68	0.76
1:B:67:LYS:H	1:B:70:HIS:HD2	1.37	0.70
1:A:116:LEU:HB3	1:A:117:PRO:HD3	1.76	0.68
1:B:267:SER:HB3	1:B:299:LEU:HG	1.75	0.68
1:A:105:MET:HE1	1:B:115:GLU:HB3	1.78	0.65
1:A:200:THR:HG23	1:A:219:TYR:CD2	2.33	0.64
1:A:64:GLU:O	1:A:85:ASP:HB2	1.99	0.62
1:A:344:LEU:O	1:A:348:ILE:HG12	1.99	0.62
1:B:97:HIS:HD2	1:B:232:ASP:O	1.83	0.62
1:A:237:ILE:HB	1:A:240:ILE:HD11	1.84	0.60
1:B:64:GLU:O	1:B:85:ASP:HB2	2.02	0.60
1:B:344:LEU:O	1:B:348:ILE:HG12	2.03	0.59
1:A:42:TYR:HB2	1:A:70:HIS:CD2	2.38	0.58
1:A:116:LEU:O	1:A:120:LYS:HB2	2.04	0.57
1:A:116:LEU:HB3	1:A:117:PRO:CD	2.34	0.57
1:B:191:GLY:HA3	1:B:300:ASN:HD22	1.71	0.56
1:B:256:LYS:HD2	1:B:288:ILE:O	2.05	0.56
1:B:266:ASP:HB3	1:B:269:PHE:CE2	2.41	0.56
1:A:275:ILE:HG23	1:A:319:LEU:HD21	1.87	0.56
1:B:191:GLY:HA3	1:B:300:ASN:ND2	2.21	0.55
1:B:62:ILE:HD12	1:B:66:LEU:HD11	1.88	0.55
1:B:116:LEU:HB2	1:B:117:PRO:HD3	1.89	0.55
1:A:331:ILE:HD11	1:A:344:LEU:HD13	1.89	0.54
1:B:282:PHE:O	1:B:286:ARG:HG3	2.07	0.54
1:A:278:LEU:HD23	1:A:319:LEU:HD23	1.88	0.54
1:B:223:ILE:HD12	1:B:228:ILE:HD12	1.91	0.53
1:A:72:ILE:O	1:A:76:ARG:HG3	2.08	0.53
1:A:105:MET:CE	1:B:115:GLU:HB3	2.38	0.53
1:B:346:ASP:O	1:B:350:GLN:HG3	2.09	0.52
1:A:182:ILE:HD12	1:A:260:ALA:HB3	1.92	0.51
1:A:261:LEU:HB2	1:A:295:ILE:HD12	1.93	0.50
1:B:269:PHE:HB3	1:B:273:LEU:HB2	1.94	0.50
1:B:90:LEU:HD22	1:B:237:ILE:HD13	1.92	0.50
1:A:313:LEU:HD13	1:A:332:PRO:HG3	1.94	0.49
1:A:97:HIS:HD2	1:A:232:ASP:O	1.96	0.49
1:A:41:PHE:O	1:A:42:TYR:HB3	2.13	0.48
1:B:44:GLN:HB3	1:B:46:ASP:OD1	2.13	0.48
1:A:242:PRO:HA	1:A:245:VAL:HG23	1.96	0.48
1:B:202:LEU:HD13	1:B:216:PRO:HD2	1.95	0.48
1:A:221:ILE:HD13	1:A:230:LEU:HD11	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:ILE:HD12	1:B:260:ALA:HB3	1.97	0.47
1:A:283:GLU:O	1:A:287:GLU:HG3	2.15	0.46
1:A:112:LEU:O	1:A:116:LEU:HB2	2.15	0.46
1:B:72:ILE:O	1:B:76:ARG:HB2	2.15	0.46
1:B:195:LEU:C	1:B:195:LEU:HD23	2.41	0.46
1:B:331:ILE:HA	1:B:332:PRO:HD3	1.83	0.46
1:B:59:THR:HG23	1:B:82:GLU:HB3	1.98	0.46
1:A:68:PRO:HB2	1:A:121:GLU:OE1	2.17	0.45
1:A:288:ILE:HG13	1:A:290:VAL:HG23	1.98	0.45
1:A:228:ILE:HD12	1:A:228:ILE:N	2.32	0.45
1:B:47:LYS:O	1:B:51:ILE:HG12	2.16	0.45
1:B:68:PRO:O	1:B:72:ILE:HG13	2.16	0.45
1:B:278:LEU:HD11	1:B:316:VAL:HG22	1.99	0.45
1:A:261:LEU:HB2	1:A:295:ILE:CD1	2.48	0.44
1:B:186:GLY:N	1:B:192:LYS:HD2	2.33	0.44
1:A:113:LYS:O	1:A:117:PRO:CD	2.56	0.44
1:B:242:PRO:HA	1:B:245:VAL:HG23	2.00	0.44
1:B:184:ILE:HD13	1:B:195:LEU:HD22	1.98	0.44
1:A:72:ILE:CG2	1:A:76:ARG:HH21	2.31	0.43
1:A:121:GLU:HG2	1:A:121:GLU:O	2.18	0.43
1:A:37:PRO:HB3	1:A:42:TYR:CD1	2.54	0.43
1:B:34:PRO:HB3	1:B:44:GLN:OE1	2.18	0.43
1:A:104:LYS:NZ	1:A:104:LYS:HB3	2.33	0.43
1:A:72:ILE:HG21	1:A:76:ARG:HH21	1.84	0.43
1:B:192:LYS:HB3	1:B:192:LYS:HE2	1.79	0.42
1:A:87:ILE:HD12	1:A:114:TYR:CE1	2.54	0.42
1:A:188:THR:HG22	1:A:188:THR:O	2.19	0.42
1:A:192:LYS:HE2	1:A:192:LYS:HB3	1.78	0.42
1:B:4:ALA:HA	1:B:59:THR:O	2.19	0.42
1:A:8:VAL:O	1:A:32:LYS:HA	2.19	0.42
1:B:13:GLU:O	1:B:17:ILE:HG13	2.20	0.42
1:B:220:ALA:HB2	1:B:229:MET:HE2	2.00	0.42
1:B:331:ILE:HD11	1:B:344:LEU:HD13	2.02	0.42
1:B:87:ILE:HD12	1:B:114:TYR:CE1	2.54	0.42
1:B:221:ILE:HA	1:B:222:PRO:HD3	1.84	0.42
1:A:223:ILE:HG21	1:A:349:TYR:HB2	2.02	0.42
1:A:221:ILE:HA	1:A:222:PRO:HD3	1.83	0.41
1:A:102:GLU:HB2	1:B:119:ILE:HD13	2.01	0.41
1:A:66:LEU:HB2	1:A:71:PHE:CE2	2.55	0.41
1:A:101:LYS:HE3	1:A:101:LYS:HB2	1.91	0.41
1:A:192:LYS:NZ	1:A:233:THR:O	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:GLU:O	1:B:96:LEU:HG	2.21	0.41
1:B:68:PRO:HB3	1:B:118:ILE:HA	2.03	0.41
1:A:331:ILE:HA	1:A:332:PRO:HD3	1.82	0.41
1:A:341:LEU:O	1:A:345:ARG:HG3	2.20	0.41
1:B:187:TYR:CE2	1:B:238:ARG:HG2	2.56	0.41
1:B:234:VAL:HA	1:B:235:PRO:HD3	1.88	0.41
1:B:283:GLU:HA	1:B:286:ARG:NH1	2.36	0.41
1:B:218:ARG:HD3	2:B:380:HOH:O	2.20	0.40
1:B:103:ALA:O	1:B:107:ILE:HG13	2.20	0.40
1:A:253:SER:O	1:A:256:LYS:HG2	2.22	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	283/364 (78%)	268 (95%)	15 (5%)	0	100	100
1	B	280/364 (77%)	268 (96%)	12 (4%)	0	100	100
All	All	563/728 (77%)	536 (95%)	27 (5%)	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	268/332 (81%)	267 (100%)	1 (0%)	84	91
1	B	260/332 (78%)	260 (100%)	0	100	100
All	All	528/664 (80%)	527 (100%)	1 (0%)	87	94

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	50	GLN
1	A	73	ASN
1	A	224	ASN
1	A	243	GLN
1	A	272	ASN
1	A	354	GLN
1	B	49	GLN
1	B	50	GLN
1	B	65	GLN
1	B	70	HIS
1	B	73	ASN
1	B	97	HIS
1	B	189	ASN
1	B	224	ASN
1	B	225	ASN
1	B	272	ASN

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 [i](#)

There are no oligosaccharides 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 [i](#)

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	293/364 (80%)	0.74	27 (9%) 14 17	16, 45, 74, 90	6 (2%)
1	B	283/364 (77%)	1.08	46 (16%) 4 5	18, 54, 81, 91	15 (5%)
All	All	576/728 (79%)	0.91	73 (12%) 8 9	16, 50, 79, 91	21 (3%)

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	43	ILE	5.3
1	B	28	THR	4.8
1	A	39	VAL	4.5
1	B	33	LEU	4.5
1	A	118	ILE	4.3
1	B	29	SER	4.1
1	B	189	ASN	4.1
1	B	45	TYR	3.7
1	B	336	LEU	3.6
1	A	46	ASP	3.6
1	A	122	THR	3.5
1	A	339	THR	3.5
1	B	305	ILE	3.5
1	A	310	TYR	3.2
1	B	78	LEU	3.2
1	B	122	THR	3.2
1	A	188	THR	3.1
1	A	308	ASP	3.0
1	B	27	VAL	3.0
1	A	270	SER	3.0
1	A	177	ASN	3.0
1	A	202	LEU	2.9
1	B	80	GLY	2.9
1	B	309	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	306	ASN	2.8
1	A	40	LYS	2.7
1	A	38	ASN	2.7
1	B	221	ILE	2.7
1	B	17	ILE	2.7
1	A	341	LEU	2.7
1	B	25	TYR	2.7
1	B	19	LEU	2.5
1	B	31	TYR	2.5
1	B	9	SER	2.5
1	B	8	VAL	2.5
1	B	179	ILE	2.5
1	B	6	LEU	2.5
1	B	3	THR	2.4
1	B	223	ILE	2.4
1	B	216	PRO	2.4
1	A	119	ILE	2.4
1	B	188	THR	2.4
1	B	274	LEU	2.4
1	B	30	ILE	2.3
1	B	215	SER	2.3
1	B	53	ASN	2.3
1	B	326	PRO	2.3
1	B	5	ALA	2.3
1	A	76	ARG	2.3
1	A	121	GLU	2.3
1	A	357	LEU	2.3
1	A	53	ASN	2.3
1	A	1	MET	2.3
1	B	18	ALA	2.3
1	B	10	LYS	2.3
1	B	76	ARG	2.2
1	B	310	TYR	2.2
1	B	79	LYS	2.2
1	B	38	ASN	2.2
1	B	34	PRO	2.2
1	A	79	LYS	2.2
1	A	101	LYS	2.2
1	A	311	LYS	2.1
1	B	249	PHE	2.1
1	B	202	LEU	2.1
1	B	15	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	2	LYS	2.0
1	B	21	GLU	2.0
1	B	299	LEU	2.0
1	B	290	VAL	2.0
1	A	49	GLN	2.0
1	B	44	GLN	2.0
1	A	304	LYS	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 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.