



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 17, 2016 – 10:03 AM EDT

PDB ID : 5CGA
Title : Structure of Hydroxyethylthiazole kinase ThiM from *Staphylococcus aureus* in complex with substrate analog 2-(1,3,5-trimethyl-1H-pyrazole-4-yl)ethanol
Authors : Kuenz, M.; Drebes, J.; Windshuegel, B.; Cang, H.; Wrenger, C.; Betzel, C.
Deposited on : 2015-07-09
Resolution : 1.87 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

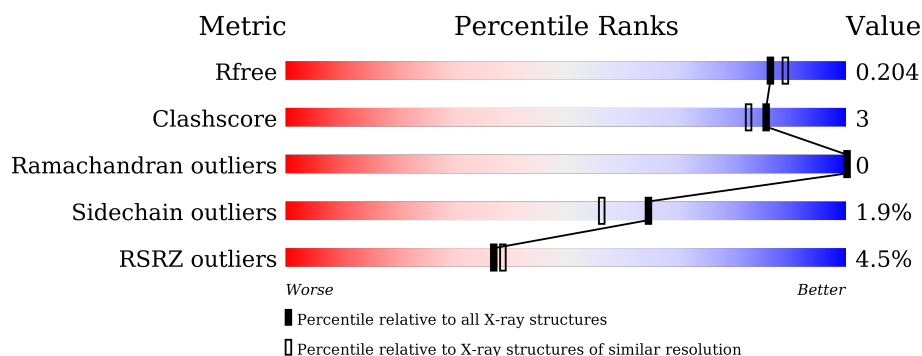
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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	6965 (1.90-1.86)
Clashscore	102246	7778 (1.90-1.86)
Ramachandran outliers	100387	7691 (1.90-1.86)
Sidechain outliers	100360	7692 (1.90-1.86)
RSRZ outliers	91569	6979 (1.90-1.86)

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	277	<div> <div>81%</div> <div>10% • 9%</div> </div>
1	B	277	<div> <div>79%</div> <div>9% • 10%</div> </div>
1	C	277	<div> <div>81%</div> <div>8% • 9%</div> </div>
1	E	277	<div> <div>4%</div> <div>85%</div> <div>6% • 8%</div> </div>
1	F	277	<div> <div>8%</div> <div>80%</div> <div>7% • 12%</div> </div>
1	H	277	<div> <div>9%</div> <div>83%</div> <div>5% 11%</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	KP6	A	301	-	-	-	X
2	KP6	B	301	-	X	-	X
2	KP6	E	301	-	-	-	X
2	KP6	F	301	-	-	-	X
2	KP6	H	301	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11784 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 Hydroxyethylthiazole kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	54	0	0
			1880	1200	309	365	6			
1	B	249	Total	C	N	O	S	34	0	0
			1862	1190	305	361	6			
1	C	251	Total	C	N	O	S	38	0	0
			1882	1199	308	369	6			
1	E	255	Total	C	N	O	S	69	0	0
			1925	1228	314	377	6			
1	F	245	Total	C	N	O	S	81	0	0
			1834	1171	301	357	5			
1	H	246	Total	C	N	O	S	34	0	0
			1845	1179	302	358	6			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	264	GLU	-	expression tag	UNP Q6GEY3
A	265	ASN	-	expression tag	UNP Q6GEY3
A	266	LEU	-	expression tag	UNP Q6GEY3
A	267	TYR	-	expression tag	UNP Q6GEY3
A	268	GLN	-	expression tag	UNP Q6GEY3
A	269	GLN	-	expression tag	UNP Q6GEY3
A	270	SER	-	expression tag	UNP Q6GEY3
A	271	GLY	-	expression tag	UNP Q6GEY3
A	272	HIS	-	expression tag	UNP Q6GEY3
A	273	HIS	-	expression tag	UNP Q6GEY3
A	274	HIS	-	expression tag	UNP Q6GEY3
A	275	HIS	-	expression tag	UNP Q6GEY3
A	276	HIS	-	expression tag	UNP Q6GEY3
A	277	HIS	-	expression tag	UNP Q6GEY3
B	264	GLU	-	expression tag	UNP Q6GEY3
B	265	ASN	-	expression tag	UNP Q6GEY3
B	266	LEU	-	expression tag	UNP Q6GEY3

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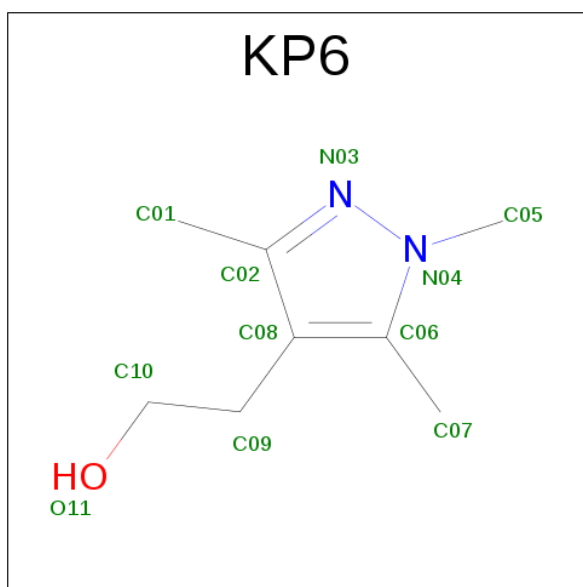
Chain	Residue	Modelled	Actual	Comment	Reference
B	267	TYR	-	expression tag	UNP Q6GEY3
B	268	GLN	-	expression tag	UNP Q6GEY3
B	269	GLN	-	expression tag	UNP Q6GEY3
B	270	SER	-	expression tag	UNP Q6GEY3
B	271	GLY	-	expression tag	UNP Q6GEY3
B	272	HIS	-	expression tag	UNP Q6GEY3
B	273	HIS	-	expression tag	UNP Q6GEY3
B	274	HIS	-	expression tag	UNP Q6GEY3
B	275	HIS	-	expression tag	UNP Q6GEY3
B	276	HIS	-	expression tag	UNP Q6GEY3
B	277	HIS	-	expression tag	UNP Q6GEY3
C	264	GLU	-	expression tag	UNP Q6GEY3
C	265	ASN	-	expression tag	UNP Q6GEY3
C	266	LEU	-	expression tag	UNP Q6GEY3
C	267	TYR	-	expression tag	UNP Q6GEY3
C	268	GLN	-	expression tag	UNP Q6GEY3
C	269	GLN	-	expression tag	UNP Q6GEY3
C	270	SER	-	expression tag	UNP Q6GEY3
C	271	GLY	-	expression tag	UNP Q6GEY3
C	272	HIS	-	expression tag	UNP Q6GEY3
C	273	HIS	-	expression tag	UNP Q6GEY3
C	274	HIS	-	expression tag	UNP Q6GEY3
C	275	HIS	-	expression tag	UNP Q6GEY3
C	276	HIS	-	expression tag	UNP Q6GEY3
C	277	HIS	-	expression tag	UNP Q6GEY3
E	264	GLU	-	expression tag	UNP Q6GEY3
E	265	ASN	-	expression tag	UNP Q6GEY3
E	266	LEU	-	expression tag	UNP Q6GEY3
E	267	TYR	-	expression tag	UNP Q6GEY3
E	268	GLN	-	expression tag	UNP Q6GEY3
E	269	GLN	-	expression tag	UNP Q6GEY3
E	270	SER	-	expression tag	UNP Q6GEY3
E	271	GLY	-	expression tag	UNP Q6GEY3
E	272	HIS	-	expression tag	UNP Q6GEY3
E	273	HIS	-	expression tag	UNP Q6GEY3
E	274	HIS	-	expression tag	UNP Q6GEY3
E	275	HIS	-	expression tag	UNP Q6GEY3
E	276	HIS	-	expression tag	UNP Q6GEY3
E	277	HIS	-	expression tag	UNP Q6GEY3
F	264	GLU	-	expression tag	UNP Q6GEY3
F	265	ASN	-	expression tag	UNP Q6GEY3
F	266	LEU	-	expression tag	UNP Q6GEY3

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Chain	Residue	Modelled	Actual	Comment	Reference
F	267	TYR	-	expression tag	UNP Q6GEY3
F	268	GLN	-	expression tag	UNP Q6GEY3
F	269	GLN	-	expression tag	UNP Q6GEY3
F	270	SER	-	expression tag	UNP Q6GEY3
F	271	GLY	-	expression tag	UNP Q6GEY3
F	272	HIS	-	expression tag	UNP Q6GEY3
F	273	HIS	-	expression tag	UNP Q6GEY3
F	274	HIS	-	expression tag	UNP Q6GEY3
F	275	HIS	-	expression tag	UNP Q6GEY3
F	276	HIS	-	expression tag	UNP Q6GEY3
F	277	HIS	-	expression tag	UNP Q6GEY3
H	264	GLU	-	expression tag	UNP Q6GEY3
H	265	ASN	-	expression tag	UNP Q6GEY3
H	266	LEU	-	expression tag	UNP Q6GEY3
H	267	TYR	-	expression tag	UNP Q6GEY3
H	268	GLN	-	expression tag	UNP Q6GEY3
H	269	GLN	-	expression tag	UNP Q6GEY3
H	270	SER	-	expression tag	UNP Q6GEY3
H	271	GLY	-	expression tag	UNP Q6GEY3
H	272	HIS	-	expression tag	UNP Q6GEY3
H	273	HIS	-	expression tag	UNP Q6GEY3
H	274	HIS	-	expression tag	UNP Q6GEY3
H	275	HIS	-	expression tag	UNP Q6GEY3
H	276	HIS	-	expression tag	UNP Q6GEY3
H	277	HIS	-	expression tag	UNP Q6GEY3

- Molecule 2 is 2-(1,3,5-trimethyl-1H-pyrazol-4-yl)ethanol (three-letter code: KP6) (formula: C₈H₁₄N₂O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	8	2	1		
2	B	1	Total	C	N	O	0	0
			11	8	2	1		
2	C	1	Total	C	N	O	0	0
			11	8	2	1		
2	E	1	Total	C	N	O	0	0
			11	8	2	1		
2	F	1	Total	C	N	O	0	0
			11	8	2	1		
2	H	1	Total	C	N	O	0	0
			11	8	2	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		

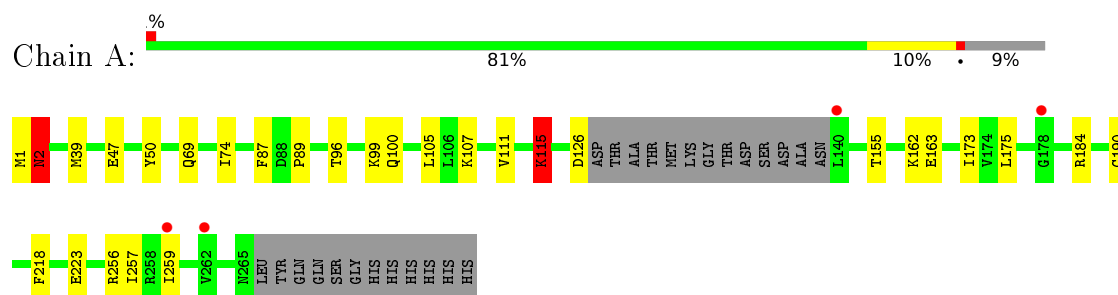
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	115	Total 115	O 115	0	0
4	B	134	Total 134	O 134	0	0
4	C	140	Total 140	O 140	0	0
4	E	49	Total 49	O 49	0	0
4	F	25	Total 25	O 25	0	0
4	H	23	Total 23	O 23	0	0

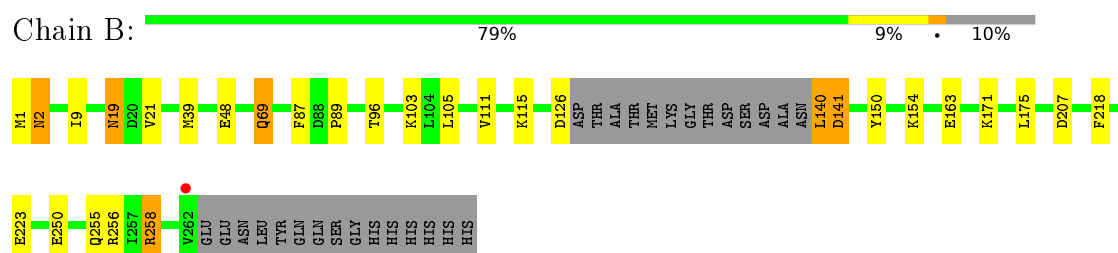
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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.

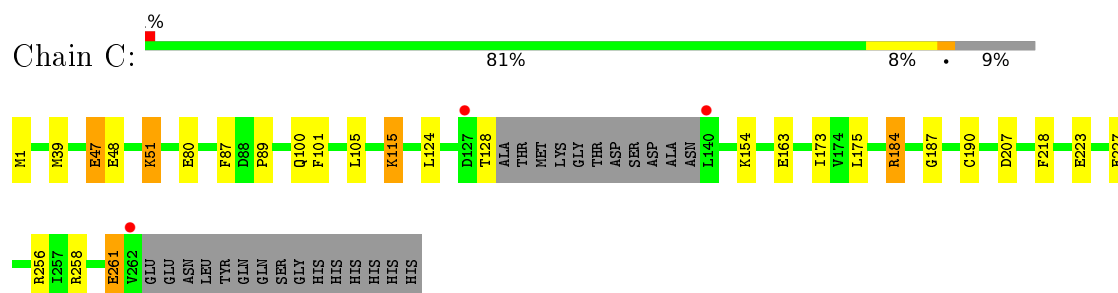
• Molecule 1: Hydroxyethylthiazole kinase



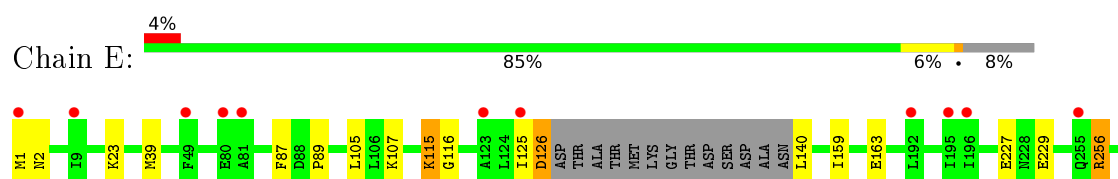
• Molecule 1: Hydroxyethylthiazole kinase

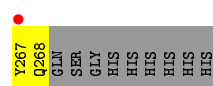


• Molecule 1: Hydroxyethylthiazole kinase

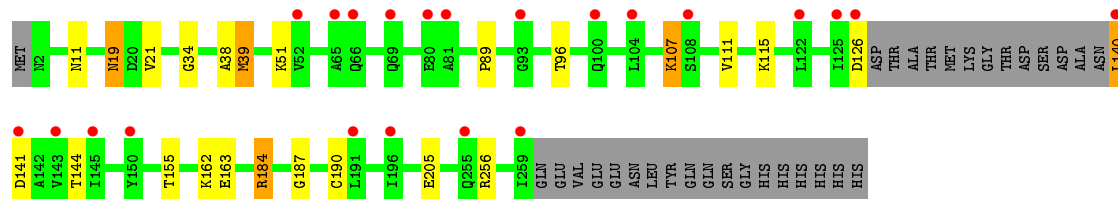
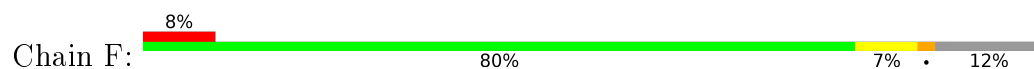


• Molecule 1: Hydroxyethylthiazole kinase

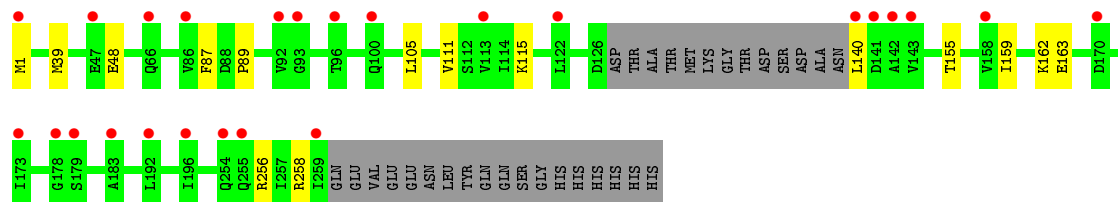
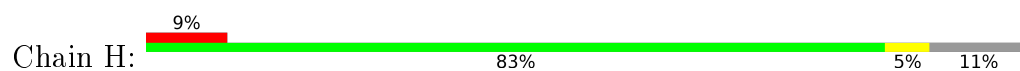




● Molecule 1: Hydroxyethylthiazole kinase



● Molecule 1: Hydroxyethylthiazole kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.03Å 62.36Å 108.32Å 92.59° 91.39° 101.28°	Depositor
Resolution (Å)	29.81 – 1.87 29.79 – 1.87	Depositor EDS
% Data completeness (in resolution range)	95.8 (29.81-1.87) 94.2 (29.79-1.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.169 , 0.199 0.176 , 0.204	Depositor DCC
R_{free} test set	6337 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.6	EDS
Estimated twinning fraction	0.025 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 125935 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11784	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, KP6

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	1.48	13/1904 (0.7%)	1.20	13/2591 (0.5%)
1	B	1.26	9/1886 (0.5%)	1.36	14/2566 (0.5%)
1	C	1.53	11/1906 (0.6%)	1.66	15/2593 (0.6%)
1	E	1.15	8/1950 (0.4%)	1.07	13/2652 (0.5%)
1	F	1.39	6/1858 (0.3%)	3.33	10/2528 (0.4%)
1	H	1.05	3/1869 (0.2%)	1.30	5/2542 (0.2%)
All	All	1.32	50/11373 (0.4%)	1.81	70/15472 (0.5%)

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	1
1	B	0	4
1	C	0	2
1	E	0	1
1	F	0	1
All	All	0	9

The worst 5 of 50 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	184	ARG	NE-CZ	-37.92	0.83	1.33
1	A	184	ARG	CZ-NH2	-33.76	0.89	1.33
1	F	126	ASP	CA-C	31.07	2.33	1.52
1	C	258	ARG	NE-CZ	-28.20	0.96	1.33
1	H	258	ARG	CZ-NH1	-27.63	0.97	1.33

The worst 5 of 70 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	184	ARG	NE-CZ-NH1	-149.18	45.71	120.30
1	H	258	ARG	NE-CZ-NH2	-48.75	95.92	120.30
1	F	126	ASP	CA-C-O	-44.99	25.61	120.10
1	C	184	ARG	NE-CZ-NH2	-43.97	98.31	120.30
1	C	184	ARG	NE-CZ-NH1	-35.03	102.78	120.30

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	ASN	Sidechain
1	B	141	ASP	Sidechain
1	B	2	ASN	Sidechain
1	B	255	GLN	Sidechain
1	B	258	ARG	Sidechain

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	1880	0	1917	12	0
1	B	1862	0	1907	9	0
1	C	1882	0	1919	10	0
1	E	1925	0	1963	6	0
1	F	1834	0	1873	24	0
1	H	1845	0	1894	6	0
2	A	11	0	0	0	0
2	B	11	0	0	0	0
2	C	11	0	0	0	0
2	E	11	0	0	5	0
2	F	11	0	0	0	0
2	H	11	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
4	A	115	0	0	0	0
4	B	134	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	140	0	0	3	1
4	E	49	0	0	0	1
4	F	25	0	0	2	0
4	H	23	0	0	1	0
All	All	11784	0	11473	62	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 3.

The worst 5 of 62 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:F:39:MET:HE2	1:F:39:MET:N	1.45	1.28
1:F:38:ALA:C	1:F:39:MET:HE2	1.80	0.99
1:F:39:MET:CE	1:F:39:MET:N	2.28	0.96
1:F:39:MET:HE2	1:F:39:MET:H	1.44	0.82
1:A:175:LEU:CD2	1:A:259:ILE:HG22	2.10	0.81

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)
4:C:426:HOH:O	4:E:413:HOH:O[1_556]	2.17	0.03

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	248/277 (90%)	245 (99%)	3 (1%)	0	100	100
1	B	245/277 (88%)	243 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	247/277 (89%)	246 (100%)	1 (0%)	0	100	100
1	E	251/277 (91%)	248 (99%)	3 (1%)	0	100	100
1	F	241/277 (87%)	239 (99%)	2 (1%)	0	100	100
1	H	242/277 (87%)	240 (99%)	2 (1%)	0	100	100
All	All	1474/1662 (89%)	1461 (99%)	13 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	198/223 (89%)	194 (98%)	4 (2%)	63	54
1	B	197/223 (88%)	191 (97%)	6 (3%)	48	36
1	C	200/223 (90%)	197 (98%)	3 (2%)	72	66
1	E	205/223 (92%)	203 (99%)	2 (1%)	82	79
1	F	194/223 (87%)	189 (97%)	5 (3%)	54	42
1	H	196/223 (88%)	193 (98%)	3 (2%)	72	66
All	All	1190/1338 (89%)	1167 (98%)	23 (2%)	65	56

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	39	MET
1	C	115	LYS
1	H	140	LEU
1	C	51	LYS
1	E	39	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	19	ASN
1	H	67	ASN
1	H	6	ASN
1	E	260	GLN
1	H	2	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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$
2	KP6	A	301	-	11,11,11	2.32	3 (27%)	12,15,15	2.75	5 (41%)
2	KP6	B	301	-	11,11,11	2.60	5 (45%)	12,15,15	4.13	7 (58%)
2	KP6	C	301	-	11,11,11	2.33	4 (36%)	12,15,15	2.97	5 (41%)
2	KP6	E	301	-	11,11,11	2.49	5 (45%)	12,15,15	1.94	5 (41%)
2	KP6	F	301	-	11,11,11	2.28	3 (27%)	12,15,15	2.38	3 (25%)
2	KP6	H	301	-	11,11,11	2.49	5 (45%)	12,15,15	2.39	5 (41%)

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	KP6	A	301	-	-	0/3/3/3	0/1/1/1
2	KP6	B	301	-	-	0/3/3/3	0/1/1/1
2	KP6	C	301	-	-	0/3/3/3	0/1/1/1
2	KP6	E	301	-	-	0/3/3/3	0/1/1/1
2	KP6	F	301	-	-	0/3/3/3	0/1/1/1
2	KP6	H	301	-	-	0/3/3/3	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	301	KP6	C06-N04	-3.02	1.33	1.37
2	B	301	KP6	N03-N04	-2.38	1.35	1.37
2	E	301	KP6	C06-N04	-2.08	1.34	1.37
2	C	301	KP6	C02-N03	2.40	1.38	1.33
2	F	301	KP6	N03-N04	2.46	1.40	1.37

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	KP6	C06-N04-N03	-5.89	107.67	111.84
2	B	301	KP6	C01-C02-C08	-4.50	119.14	129.69
2	H	301	KP6	C01-C02-C08	-3.01	122.64	129.69
2	A	301	KP6	C01-C02-C08	-2.92	122.84	129.69
2	C	301	KP6	C06-N04-N03	-2.37	110.16	111.84

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	301	KP6	5	0

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	252/277 (90%)	-0.18	4 (1%) 74 77	14, 27, 44, 62	26 (10%)
1	B	249/277 (89%)	-0.33	1 (0%) 93 93	13, 23, 37, 60	19 (7%)
1	C	251/277 (90%)	-0.25	3 (1%) 81 83	12, 24, 39, 74	24 (9%)
1	E	255/277 (92%)	0.10	12 (4%) 35 37	25, 40, 64, 75	25 (9%)
1	F	245/277 (88%)	0.52	22 (8%) 12 12	27, 51, 76, 103	22 (8%)
1	H	246/277 (88%)	0.46	25 (10%) 9 9	27, 48, 68, 89	11 (4%)
All	All	1498/1662 (90%)	0.05	67 (4%) 37 39	12, 34, 66, 103	127 (8%)

The worst 5 of 67 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	140	LEU	8.8
1	F	122	LEU	6.1
1	F	52	VAL	5.7
1	H	140	LEU	5.0
1	H	259	ILE	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 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(Å ²)	Q<0.9
2	KP6	B	301	11/11	0.92	0.21	5.47	22,30,36,38	2
2	KP6	E	301	11/11	0.89	0.25	5.21	39,40,45,48	5
2	KP6	A	301	11/11	0.90	0.17	5.10	24,40,45,48	1
2	KP6	F	301	11/11	0.93	0.24	4.99	32,38,47,48	4
2	KP6	H	301	11/11	0.91	0.20	3.30	35,42,52,52	3
2	KP6	C	301	11/11	0.93	0.13	1.94	24,31,37,39	2
3	MG	A	302	1/1	0.75	0.18	1.80	46,46,46,46	1
3	MG	E	302	1/1	0.86	0.16	0.40	47,47,47,47	1
3	MG	C	302	1/1	0.81	0.10	-0.22	43,43,43,43	1
3	MG	B	302	1/1	0.93	0.09	-0.38	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.