



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:07 AM UTC

PDB ID : 5RG6 / pdb_00005rg6
Title : Crystal Structure of Kemp Eliminase HG3.7 in unbound state, 277K
Authors : Broom, A.; Rakotoharisoa, R.V.; Thompson, M.C.; Fraser, J.S.; Chica, R.A.
Deposited on : 2020-03-19
Resolution : 1.35 Å(reported)

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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
Mogul : 2022.3.0, CSD as543be (2022)
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

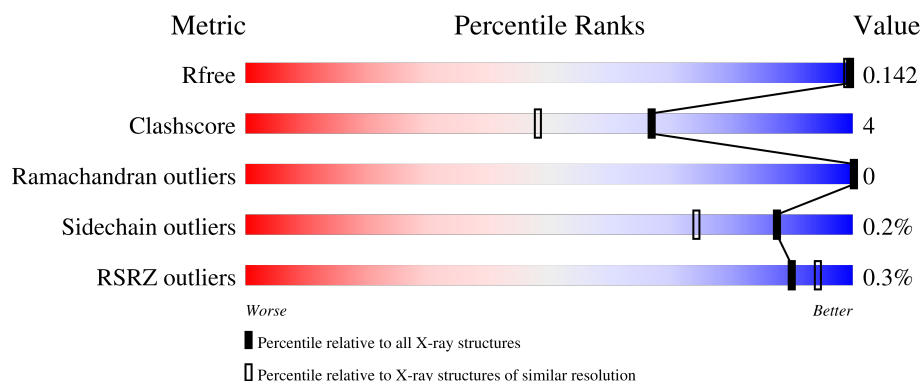
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.35 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	1216 (1.36-1.36)
Clashscore	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)
RSRZ outliers	180081	1214 (1.36-1.36)

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	318	 86% 8% 6%
1	B	318	 87% 8% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11863 atoms, of which 5634 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 Kemp Eliminate HG3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	300	Total	C	H	N	O	S	0	98	0
			5563	1758	2774	493	524	14			
1	B	300	Total	C	H	N	O	S	1	113	0
			5732	1809	2860	510	539	14			

There are 66 discrepancies between the modelled and reference sequences:

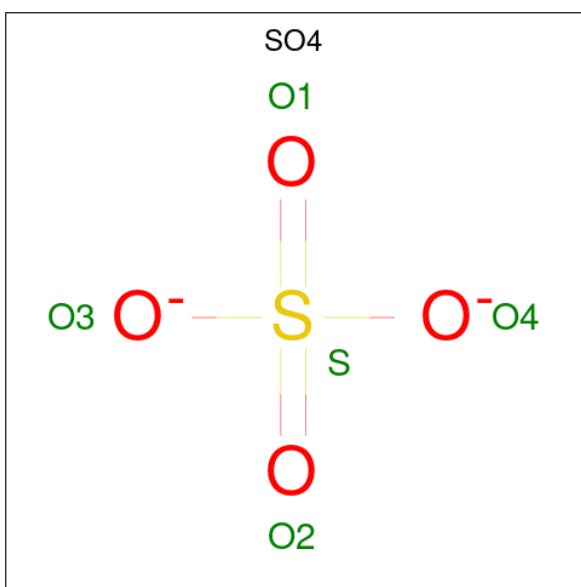
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P23360
A	0	ALA	-	expression tag	UNP P23360
A	1	GLU	-	expression tag	UNP P23360
A	6	ILE	VAL	engineered mutation	UNP P23360
A	37	LYS	GLN	engineered mutation	UNP P23360
A	42	MET	GLN	engineered mutation	UNP P23360
A	44	TRP	THR	engineered mutation	UNP P23360
A	50	GLN	LYS	engineered mutation	UNP P23360
A	81	GLY	ARG	engineered mutation	UNP P23360
A	83	GLY	HIS	engineered mutation	UNP P23360
A	84	CYS	THR	engineered mutation	UNP P23360
A	89	ARG	SER	engineered mutation	UNP P23360
A	90	HIS	GLN	engineered mutation	UNP P23360
A	125	ASN	ALA	engineered mutation	UNP P23360
A	130	GLY	ASN	engineered mutation	UNP P23360
A	172	MET	ASN	engineered mutation	UNP P23360
A	234	SER	ALA	engineered mutation	UNP P23360
A	236	LEU	THR	engineered mutation	UNP P23360
A	237	MET	GLU	engineered mutation	UNP P23360
A	267	PHE	TRP	engineered mutation	UNP P23360
A	304	GLY	-	expression tag	UNP P23360
A	305	SER	-	expression tag	UNP P23360
A	306	ILE	-	expression tag	UNP P23360
A	307	GLU	-	expression tag	UNP P23360
A	308	GLY	-	expression tag	UNP P23360

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Chain	Residue	Modelled	Actual	Comment	Reference
A	309	ARG	-	expression tag	UNP P23360
A	310	GLY	-	expression tag	UNP P23360
A	311	HIS	-	expression tag	UNP P23360
A	312	HIS	-	expression tag	UNP P23360
A	313	HIS	-	expression tag	UNP P23360
A	314	HIS	-	expression tag	UNP P23360
A	315	HIS	-	expression tag	UNP P23360
A	316	HIS	-	expression tag	UNP P23360
B	-1	MET	-	initiating methionine	UNP P23360
B	0	ALA	-	expression tag	UNP P23360
B	1	GLU	-	expression tag	UNP P23360
B	6	ILE	VAL	engineered mutation	UNP P23360
B	37	LYS	GLN	engineered mutation	UNP P23360
B	42	MET	GLN	engineered mutation	UNP P23360
B	44	TRP	THR	engineered mutation	UNP P23360
B	50	GLN	LYS	engineered mutation	UNP P23360
B	81	GLY	ARG	engineered mutation	UNP P23360
B	83	GLY	HIS	engineered mutation	UNP P23360
B	84	CYS	THR	engineered mutation	UNP P23360
B	89	ARG	SER	engineered mutation	UNP P23360
B	90	HIS	GLN	engineered mutation	UNP P23360
B	125	ASN	ALA	engineered mutation	UNP P23360
B	130	GLY	ASN	engineered mutation	UNP P23360
B	172	MET	ASN	engineered mutation	UNP P23360
B	234	SER	ALA	engineered mutation	UNP P23360
B	236	LEU	THR	engineered mutation	UNP P23360
B	237	MET	GLU	engineered mutation	UNP P23360
B	267	PHE	TRP	engineered mutation	UNP P23360
B	304	GLY	-	expression tag	UNP P23360
B	305	SER	-	expression tag	UNP P23360
B	306	ILE	-	expression tag	UNP P23360
B	307	GLU	-	expression tag	UNP P23360
B	308	GLY	-	expression tag	UNP P23360
B	309	ARG	-	expression tag	UNP P23360
B	310	GLY	-	expression tag	UNP P23360
B	311	HIS	-	expression tag	UNP P23360
B	312	HIS	-	expression tag	UNP P23360
B	313	HIS	-	expression tag	UNP P23360
B	314	HIS	-	expression tag	UNP P23360
B	315	HIS	-	expression tag	UNP P23360
B	316	HIS	-	expression tag	UNP P23360

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

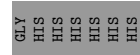


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	268	Total	O	0	9
			270	270		
3	B	270	Total	O	0	13
			273	273		

- Molecule 1: Kemp Eliminase HG3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.20Å 79.85Å 98.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.34 – 1.35 60.34 – 1.35	Depositor EDS
% Data completeness (in resolution range)	98.9 (60.34-1.35) 98.9 (60.34-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 1.35Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.126 , 0.141 0.128 , 0.142	Depositor DCC
R_{free} test set	6660 reflections (2.79%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	11863	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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	0.36	0/3188	0.55	0/4350
1	B	0.34	0/3302	0.54	0/4503
All	All	0.35	0/6490	0.54	0/8853

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	2789	2774	2435	21	0
1	B	2872	2860	2477	20	0
2	A	10	0	0	0	0
2	B	15	0	0	1	0
3	A	270	0	0	6	1
3	B	273	0	0	9	1
All	All	6229	5634	4912	42	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 4.

All (42) 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:207[B]:GLN:NE2	1:A:237[B]:MET:SD	2.25	1.10
1:B:237[B]:MET:SD	3:B:537:HOH:O	2.25	0.93
1:B:236[B]:LEU:HD21	3:B:519:HOH:O	1.71	0.91
1:A:236[B]:LEU:HD21	3:A:537:HOH:O	1.75	0.86
1:B:125[A]:ASN:ND2	3:B:501:HOH:O	1.97	0.85
1:A:237[B]:MET:SD	3:A:599:HOH:O	2.36	0.83
1:A:47[B]:ASN:OD1	3:A:501:HOH:O	2.08	0.71
1:B:100:ASP:OD2	1:B:103[B]:THR:HG23	1.90	0.70
1:B:17:TYR:OH	3:B:501:HOH:O	2.09	0.68
1:B:44[B]:TRP:NE1	1:B:265[B]:THR:HG21	2.11	0.65
1:B:26[B]:ARG:NE	1:B:46[B]:GLU:OE1	2.26	0.64
1:A:275[B]:TRP:CH2	1:A:276[B]:ARG:HD2	2.35	0.62
1:A:117[B]:THR:HG22	3:A:543:HOH:O	1.98	0.61
1:B:89[B]:ARG:O	3:B:502:HOH:O	2.15	0.61
1:B:84[B]:CYS:SG	1:B:87[B]:TRP:HB2	2.41	0.61
1:B:113[B]:THR:HG22	3:B:634:HOH:O	2.01	0.59
1:A:21:ALA:HB3	1:A:267[B]:PHE:O	2.04	0.57
1:A:84[B]:CYS:SG	1:A:87[B]:TRP:HB2	2.45	0.56
1:B:113[B]:THR:HG21	3:B:517:HOH:O	2.06	0.56
1:A:47[B]:ASN:OD1	1:A:50[B]:GLN:OE1	2.26	0.52
1:B:89[A]:ARG:NH2	2:B:403:SO4:O3	2.42	0.49
1:B:21:ALA:HB3	1:B:267[B]:PHE:O	2.12	0.48
1:B:195[A]:ARG:HD2	1:B:230:THR:HG22	1.96	0.48
1:A:54:THR:O	1:A:61[A]:PHE:HA	2.13	0.47
1:B:299[B]:GLN:HG3	3:B:550:HOH:O	2.15	0.47
1:A:207[B]:GLN:OE1	1:A:236[B]:LEU:HB2	2.16	0.46
1:B:109[B]:LYS:O	1:B:113[B]:THR:HG23	2.16	0.45
1:A:195[B]:ARG:HD2	1:A:230:THR:HG22	2.00	0.44
1:A:236[A]:LEU:HG	1:A:237[A]:MET:HG2	2.00	0.44
1:A:275[B]:TRP:CH2	1:A:276[B]:ARG:CD	3.01	0.44
1:A:275[B]:TRP:CZ2	1:A:276[B]:ARG:HD3	2.52	0.44
1:B:54:THR:O	1:B:61:PHE:HA	2.18	0.43
1:A:47[A]:ASN:ND2	3:A:515[A]:HOH:O	2.52	0.42
1:B:44[B]:TRP:HE1	1:B:265[B]:THR:HG21	1.81	0.42
1:B:195[B]:ARG:HD2	1:B:230:THR:HG22	2.01	0.42
1:A:203:GLY:HA2	1:A:232:GLU:O	2.19	0.42
1:A:235[B]:ILE:HG13	1:A:255[B]:CYS:SG	2.60	0.42
1:A:53:ALA:O	1:A:62[A]:ASN:HB2	2.21	0.41
1:A:259[A]:GLN:HG2	3:A:665:HOH:O	2.21	0.41
1:A:44[B]:TRP:CD1	1:A:45:PRO:HD2	2.56	0.40

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)
3:A:573:HOH:O	3:B:739:HOH:O[1_655]	2.14	0.06

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	397/318 (125%)	390 (98%)	7 (2%)	0	100	100
1	B	412/318 (130%)	404 (98%)	8 (2%)	0	100	100
All	All	809/636 (127%)	794 (98%)	15 (2%)	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	334/256 (130%)	334 (100%)	0	100	100
1	B	345/256 (135%)	344 (100%)	1 (0%)	86	72
All	All	679/512 (133%)	678 (100%)	1 (0%)	87	75

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

Mol	Chain	Res	Type
1	B	62	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	GLN
1	A	189	ASN
1	B	134	ASN
1	B	287	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	SO4	B	403	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	B	402	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	B	401	-	4,4,4	0.24	0	6,6,6	0.11	0
2	SO4	A	402	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	A	401	-	4,4,4	0.25	0	6,6,6	0.08	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	403	SO4	1	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, 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	300/318 (94%)	-0.53	1 (0%) 90 94	5, 14, 25, 54	63 (21%)
1	B	300/318 (94%)	-0.39	1 (0%) 90 94	5, 15, 28, 47	72 (24%)
All	All	600/636 (94%)	-0.46	2 (0%) 90 94	5, 15, 27, 54	135 (22%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	303	GLN	3.6
1	B	303	GLN	2.6

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](#)

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
2	SO4	A	402	5/5	0.62	0.17	90,99,101,103	5
2	SO4	B	403	5/5	0.65	0.16	93,104,107,111	5
2	SO4	B	401	5/5	0.72	0.16	78,99,100,107	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	402	5/5	0.79	0.13	75,77,79,80	5
2	SO4	A	401	5/5	0.81	0.15	55,74,78,90	0

6.5 Other polymers [i](#)

There are no such residues in this entry.