



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:39 PM UTC

PDB ID : 5IDI / pdb_00005idi
Title : Structure of beta glucosidase 1A from Thermotoga neapolitana, mutant E349A
Authors : Kulkarni, T.; Nordberg Karlsson, E.; Logan, D.T.
Deposited on : 2016-02-24
Resolution : 1.90 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
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

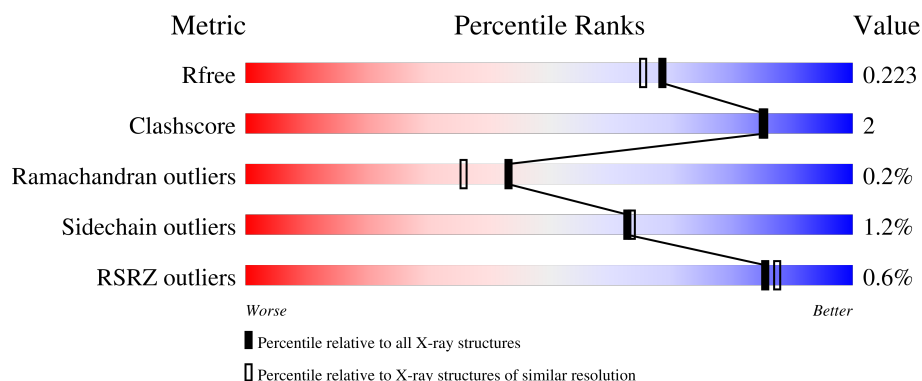
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	452	% 82% 15% ..
1	B	452	83% 15% .

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	ACT	A	501	-	-	X	-
2	ACT	A	502	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7747 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 1,4-beta-D-glucan glucohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3646	2364	616	658	8			
1	B	444	Total	C	N	O	S	0	1	0
			3660	2372	617	663	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	165	LEU	PRO	conflict	UNP B9K7M5
A	349	GLY	GLU	engineered mutation	UNP B9K7M5
A	445	LEU	-	expression tag	UNP B9K7M5
A	446	GLU	-	expression tag	UNP B9K7M5
A	447	HIS	-	expression tag	UNP B9K7M5
A	448	HIS	-	expression tag	UNP B9K7M5
A	449	HIS	-	expression tag	UNP B9K7M5
A	450	HIS	-	expression tag	UNP B9K7M5
A	451	HIS	-	expression tag	UNP B9K7M5
A	452	HIS	-	expression tag	UNP B9K7M5
B	165	LEU	PRO	conflict	UNP B9K7M5
B	349	GLY	GLU	engineered mutation	UNP B9K7M5
B	445	LEU	-	expression tag	UNP B9K7M5
B	446	GLU	-	expression tag	UNP B9K7M5
B	447	HIS	-	expression tag	UNP B9K7M5
B	448	HIS	-	expression tag	UNP B9K7M5
B	449	HIS	-	expression tag	UNP B9K7M5
B	450	HIS	-	expression tag	UNP B9K7M5
B	451	HIS	-	expression tag	UNP B9K7M5
B	452	HIS	-	expression tag	UNP B9K7M5

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

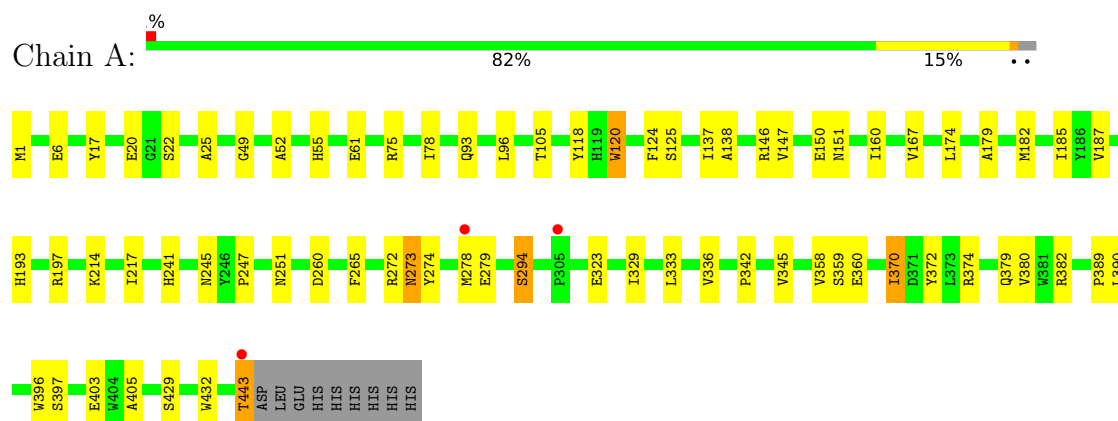
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	193	Total	O	0	0
			193	193		
3	B	236	Total	O	0	0
			236	236		

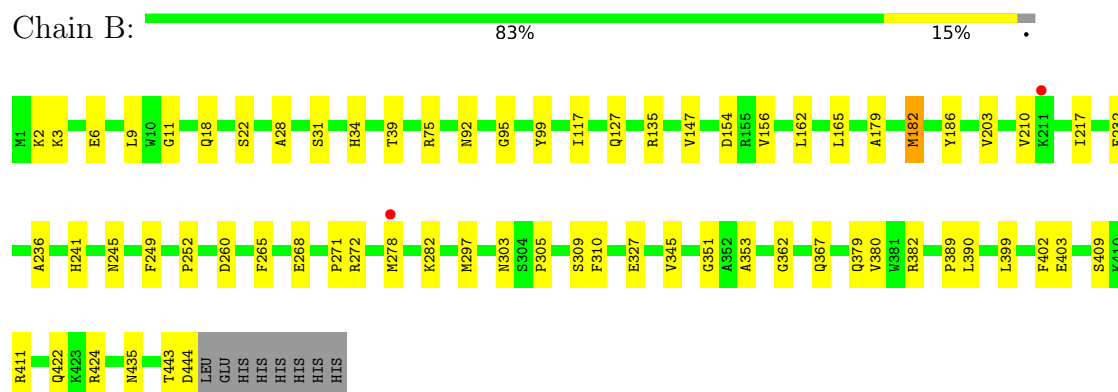
3 Residue-property plots

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: 1,4-beta-D-glucan glucohydrolase



- Molecule 1: 1,4-beta-D-glucan glucohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.28Å 98.73Å 154.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 30.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-1.90) 99.4 (30.00-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	11.63 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.169 , 0.215 0.179 , 0.223	Depositor DCC
R_{free} test set	4096 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.3	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	0.96	EDS
Total number of atoms	7747	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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](#)

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

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	1.68	35/3759 (0.9%)	1.34	18/5104 (0.4%)
1	B	1.75	46/3776 (1.2%)	1.32	7/5127 (0.1%)
All	All	1.71	81/7535 (1.1%)	1.33	25/10231 (0.2%)

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

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	22	SER	C-O	12.57	1.29	1.23
1	A	93	GLN	CA-CB	10.82	1.70	1.53
1	B	389	PRO	CA-C	7.34	1.59	1.52
1	B	22	SER	N-CA	7.14	1.52	1.46
1	A	389	PRO	CA-C	7.12	1.59	1.52
1	B	210	VAL	C-O	-7.08	1.16	1.24
1	B	351	GLY	N-CA	7.00	1.52	1.45
1	B	99	TYR	N-CA	7.00	1.54	1.46
1	A	197	ARG	CZ-NH1	6.91	1.42	1.32
1	A	151	ASN	C-O	6.88	1.33	1.24
1	B	367	GLN	CA-C	6.83	1.61	1.52
1	A	25	ALA	C-O	6.72	1.32	1.23
1	B	147	VAL	C-O	6.63	1.31	1.24
1	A	150	GLU	C-O	-6.54	1.15	1.24
1	A	96	LEU	CA-C	6.50	1.61	1.52
1	B	203	VAL	CA-C	6.44	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	403	GLU	CA-C	6.42	1.60	1.53
1	B	6	GLU	C-O	-6.40	1.16	1.23
1	B	11	GLY	N-CA	6.27	1.51	1.45
1	A	251	ASN	N-CA	6.26	1.51	1.46
1	A	147	VAL	N-CA	-6.23	1.39	1.46
1	A	6	GLU	C-O	-6.20	1.16	1.23
1	B	217	ILE	CA-CB	6.16	1.61	1.55
1	B	309	SER	CB-OG	-6.16	1.29	1.42
1	B	282	LYS	CA-CB	6.15	1.60	1.53
1	B	443	THR	CA-C	6.12	1.60	1.52
1	B	424	ARG	CA-C	6.09	1.60	1.52
1	B	34	HIS	C-O	6.05	1.31	1.24
1	B	305	PRO	C-O	5.91	1.31	1.23
1	A	160	ILE	C-O	5.88	1.30	1.24
1	B	422	GLN	N-CA	5.79	1.53	1.46
1	B	345	VAL	C-O	-5.77	1.18	1.24
1	A	185	ILE	C-O	5.71	1.31	1.24
1	A	333	LEU	N-CA	-5.71	1.39	1.46
1	B	95	GLY	N-CA	-5.69	1.38	1.45
1	B	444	ASP	CB-CG	5.69	1.66	1.52
1	B	278	MET	N-CA	5.63	1.53	1.46
1	A	17	TYR	N-CA	5.61	1.53	1.46
1	A	52	ALA	C-O	-5.58	1.19	1.24
1	B	186	TYR	C-O	5.52	1.30	1.24
1	A	405	ALA	CA-C	5.51	1.60	1.52
1	B	271	PRO	CA-C	-5.48	1.45	1.52
1	B	156	VAL	CA-C	5.47	1.59	1.52
1	B	179	ALA	CA-C	5.47	1.58	1.53
1	B	327	GLU	CD-OE1	5.47	1.35	1.25
1	B	232	GLU	CD-OE1	5.44	1.35	1.25
1	B	310	PHE	C-O	-5.44	1.17	1.23
1	A	179	ALA	N-CA	5.44	1.52	1.45
1	A	138	ALA	CA-C	5.42	1.59	1.52
1	B	28	ALA	CA-C	5.41	1.59	1.52
1	A	120	TRP	C-O	5.36	1.30	1.23
1	A	278	MET	N-CA	5.32	1.52	1.46
1	B	154	ASP	CA-C	5.30	1.59	1.52
1	A	197	ARG	CZ-NH2	-5.27	1.26	1.33
1	A	217	ILE	CA-CB	5.27	1.60	1.55
1	B	147	VAL	CA-C	5.26	1.59	1.52
1	A	49	GLY	C-O	5.24	1.30	1.23
1	B	117	ILE	CA-CB	5.22	1.60	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	92	ASN	N-CA	-5.21	1.40	1.46
1	B	409	SER	CB-OG	5.21	1.52	1.42
1	B	31	SER	C-O	5.20	1.30	1.23
1	A	22	SER	CA-CB	-5.20	1.48	1.52
1	A	61	GLU	N-CA	5.20	1.52	1.46
1	A	265	PHE	N-CA	-5.19	1.40	1.46
1	A	329	ILE	N-CA	5.16	1.53	1.46
1	A	247	PRO	CA-CB	5.16	1.61	1.53
1	B	399	LEU	N-CA	-5.16	1.39	1.45
1	B	327	GLU	CA-C	5.15	1.59	1.52
1	B	265	PHE	C-O	-5.13	1.17	1.24
1	A	174	LEU	N-CA	5.12	1.52	1.46
1	A	345	VAL	N-CA	5.10	1.52	1.46
1	B	127	GLN	CD-OE1	5.10	1.33	1.23
1	B	236	ALA	CA-C	5.10	1.59	1.52
1	B	327	GLU	CG-CD	5.10	1.64	1.52
1	B	444	ASP	N-CA	5.09	1.55	1.46
1	B	39	THR	CA-C	5.05	1.58	1.52
1	A	372	TYR	CA-C	5.04	1.59	1.52
1	A	374	ARG	CZ-NH1	5.04	1.39	1.32
1	A	167	VAL	N-CA	-5.04	1.40	1.46
1	A	358	VAL	C-O	-5.03	1.18	1.24
1	A	342	PRO	N-CA	5.00	1.53	1.47

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	443	THR	CA-CB-OG1	-7.95	97.68	109.60
1	B	75	ARG	NE-CZ-NH1	-6.57	114.93	121.50
1	A	197	ARG	NE-CZ-NH2	-6.48	113.36	119.20
1	B	260	ASP	N-CA-C	6.27	118.92	111.33
1	A	124	PHE	N-CA-C	6.22	118.85	111.33
1	A	260	ASP	N-CA-C	6.19	118.03	111.28
1	A	146	ARG	NE-CZ-NH2	-6.12	113.69	119.20
1	A	75	ARG	NE-CZ-NH1	-6.09	115.41	121.50
1	A	125	SER	CA-CB-OG	-5.98	99.15	111.10
1	A	105	THR	N-CA-C	-5.78	104.58	111.69
1	A	294	SER	N-CA-C	5.78	117.57	108.96
1	B	435	ASN	N-CA-C	5.76	117.63	111.36
1	A	75	ARG	NE-CZ-NH2	5.70	124.33	119.20
1	B	75	ARG	NE-CZ-NH2	5.64	124.27	119.20
1	A	273	ASN	N-CA-C	5.53	119.31	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	9	LEU	N-CA-C	5.51	117.72	108.96
1	A	374	ARG	NE-CZ-NH2	-5.50	114.25	119.20
1	A	443	THR	CA-CB-CG2	5.45	119.77	110.50
1	A	336	VAL	N-CA-C	-5.45	105.22	110.72
1	A	370	ILE	CB-CA-C	-5.37	104.89	112.14
1	B	182	MET	N-CA-C	5.31	118.66	110.42
1	B	303	ASN	CB-CA-C	-5.30	102.85	111.06
1	A	146	ARG	CD-NE-CZ	5.25	131.76	124.40
1	A	374	ARG	CG-CD-NE	-5.20	100.57	112.00
1	A	151	ASN	N-CA-C	5.01	119.47	112.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	359	SER	Peptide

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	3646	0	3504	18	0
1	B	3660	0	3514	10	0
2	A	8	0	6	5	0
2	B	4	0	3	1	0
3	A	193	0	0	4	0
3	B	236	0	0	3	0
All	All	7747	0	7027	31	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 2.

All (31) 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:370:ILE:HD11	1:A:429:SER:HA	1.52	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:501:ACT:H1	3:B:665:HOH:O	1.73	0.86
1:B:379:GLN:HE22	1:B:382:ARG:HH11	1.24	0.84
1:A:370:ILE:CD1	1:A:429:SER:HA	2.15	0.76
2:A:501:ACT:H1	3:A:669:HOH:O	1.85	0.75
1:A:379:GLN:HE22	1:A:382:ARG:HH11	1.36	0.73
1:B:380:VAL:HG13	1:B:390:LEU:CD2	2.29	0.62
1:A:182:MET:HG2	1:B:182:MET:HG2	1.82	0.62
1:A:294:SER:OG	2:A:502:ACT:H2	2.05	0.57
2:A:502:ACT:H1	3:A:769:HOH:O	2.13	0.48
1:B:162:LEU:HB2	1:B:165:LEU:HD11	1.97	0.47
1:A:193:HIS:CE1	1:A:274:TYR:HB2	2.50	0.46
1:A:182:MET:HE1	1:A:187:VAL:HG21	1.98	0.46
1:B:241:HIS:O	1:B:245:ASN:HB2	2.14	0.46
1:A:403:GLU:OE1	2:A:501:ACT:OXT	2.34	0.45
1:A:214:LYS:HE3	3:A:731:HOH:O	2.17	0.45
1:B:268:GLU:HB2	3:B:796:HOH:O	2.18	0.44
1:A:272:ARG:HA	1:A:272:ARG:HD3	1.93	0.42
2:A:501:ACT:H2	3:A:667:HOH:O	2.18	0.42
1:B:18:GLN:HA	1:B:402:PHE:HB3	2.02	0.42
1:A:118:TYR:CZ	1:A:120:TRP:HA	2.55	0.42
1:A:294:SER:HB2	1:A:323:GLU:O	2.20	0.41
1:B:353:ALA:HA	1:B:411:ARG:O	2.20	0.41
1:A:272:ARG:O	1:A:273:ASN:HB2	2.21	0.41
1:A:370:ILE:HD12	1:A:432:TRP:CB	2.50	0.41
1:A:20:GLU:HA	1:A:55:HIS:HB3	2.03	0.41
1:A:241:HIS:O	1:A:245:ASN:HB2	2.21	0.41
1:A:380:VAL:HG13	1:A:390:LEU:CD2	2.51	0.41
1:A:396:TRP:HA	1:A:397:SER:HA	1.91	0.40
1:B:249:PHE:C	1:B:252:PRO:HD2	2.46	0.40
1:B:297:MET:HE3	3:B:647:HOH:O	2.21	0.40

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	441/452 (98%)	429 (97%)	11 (2%)	1 (0%)	43	36
1	B	443/452 (98%)	432 (98%)	10 (2%)	1 (0%)	43	36
All	All	884/904 (98%)	861 (97%)	21 (2%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	360	GLU
1	B	362	GLY

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	380/389 (98%)	375 (99%)	5 (1%)	61	61
1	B	382/389 (98%)	378 (99%)	4 (1%)	68	70
All	All	762/778 (98%)	753 (99%)	9 (1%)	63	63

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	78	ILE
1	A	137	ILE
1	A	279	GLU
1	A	443	THR
1	B	2	LYS
1	B	3	LYS
1	B	135	ARG
1	B	272	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	57	ASN
1	A	283	GLN
1	A	350	ASN
1	A	379	GLN
1	A	385	GLN
1	A	439	ASN
1	B	57	ASN
1	B	283	GLN
1	B	379	GLN
1	B	385	GLN
1	B	439	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	ACT	B	501	-	3,3,3	1.52	1 (33%)	3,3,3	1.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ACT	A	501	-	3,3,3	1.75	1 (33%)	3,3,3	1.40	0
2	ACT	A	502	-	3,3,3	0.54	0	3,3,3	2.27	2 (66%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ACT	O-C	2.62	1.33	1.22
2	B	501	ACT	O-C	2.10	1.31	1.22

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	502	ACT	OXT-C-CH3	2.87	127.09	115.05
2	A	502	ACT	O-C-CH3	-2.69	111.50	122.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	ACT	1	0
2	A	501	ACT	3	0
2	A	502	ACT	2	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	443/452 (98%)	-0.30	3 (0%) 84 86	20, 26, 41, 68	0
1	B	444/452 (98%)	-0.39	2 (0%) 87 89	18, 24, 40, 56	1 (0%)
All	All	887/904 (98%)	-0.34	5 (0%) 85 87	18, 25, 41, 68	1 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	278	MET	2.5
1	B	278	MET	2.3
1	B	211	LYS	2.3
1	A	443	THR	2.2
1	A	305	PRO	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](#)

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($\text{\AA}^2$)	Q<0.9
2	ACT	A	502	4/4	0.83	0.14	37,38,44,48	0
2	ACT	B	501	4/4	0.88	0.11	19,23,24,34	0
2	ACT	A	501	4/4	0.91	0.10	23,25,26,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.