



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:52 AM UTC

PDB ID : 2GZX / pdb_00002gzx
Title : Crystal Structure of the TatD deoxyribonuclease MW0446 from *Staphylococcus aureus*. Northeast Structural Genomics Consortium Target ZR237.
Authors : Vorobiev, S.M.; Neely, H.; Seetharaman, J.; Wang, D.; Fang, Y.; Xiao, R.; Acton, T.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2006-05-12
Resolution : 2.20 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

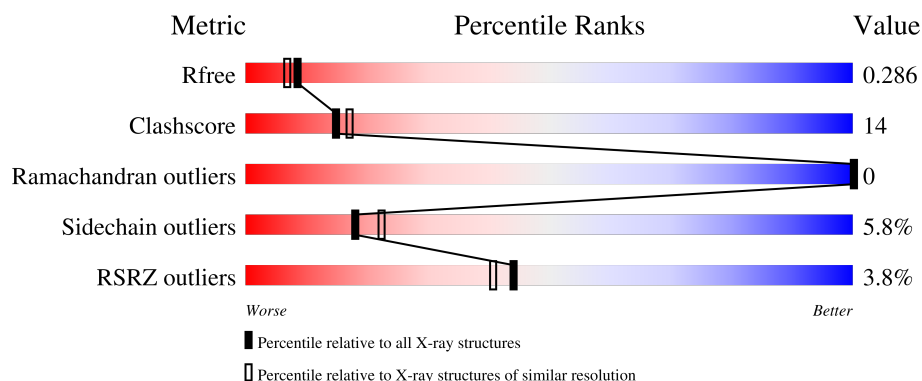
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	265	
1	B	265	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4104 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 putative TatD related DNase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	253	Total	C	N	O	S	Se	0	0	0
			2015	1283	343	382	2	5			
1	B	253	Total	C	N	O	S	Se	0	0	0
			2004	1276	341	380	2	5			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP Q1Y6Z0
A	33	MSE	MET	modified residue	UNP Q1Y6Z0
A	47	MSE	MET	modified residue	UNP Q1Y6Z0
A	93	MSE	MET	modified residue	UNP Q1Y6Z0
A	152	MSE	MET	modified residue	UNP Q1Y6Z0
A	196	MSE	MET	modified residue	UNP Q1Y6Z0
A	258	LEU	-	expression tag	UNP Q1Y6Z0
A	259	GLU	-	expression tag	UNP Q1Y6Z0
A	260	HIS	-	expression tag	UNP Q1Y6Z0
A	261	HIS	-	expression tag	UNP Q1Y6Z0
A	262	HIS	-	expression tag	UNP Q1Y6Z0
A	263	HIS	-	expression tag	UNP Q1Y6Z0
A	264	HIS	-	expression tag	UNP Q1Y6Z0
A	265	HIS	-	expression tag	UNP Q1Y6Z0
B	1	MSE	MET	modified residue	UNP Q1Y6Z0
B	33	MSE	MET	modified residue	UNP Q1Y6Z0
B	47	MSE	MET	modified residue	UNP Q1Y6Z0
B	93	MSE	MET	modified residue	UNP Q1Y6Z0
B	152	MSE	MET	modified residue	UNP Q1Y6Z0
B	196	MSE	MET	modified residue	UNP Q1Y6Z0
B	258	LEU	-	expression tag	UNP Q1Y6Z0
B	259	GLU	-	expression tag	UNP Q1Y6Z0
B	260	HIS	-	expression tag	UNP Q1Y6Z0
B	261	HIS	-	expression tag	UNP Q1Y6Z0
B	262	HIS	-	expression tag	UNP Q1Y6Z0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	263	HIS	-	expression tag	UNP Q1Y6Z0
B	264	HIS	-	expression tag	UNP Q1Y6Z0
B	265	HIS	-	expression tag	UNP Q1Y6Z0

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total 2	Ni 2	0	0
2	B	2	Total 2	Ni 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total 41	O 41	0	0
3	B	40	Total 40	O 40	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.68Å 77.32Å 76.13Å 90.00° 98.01° 90.00°	Depositor
Resolution (Å)	26.79 – 2.20 26.79 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (26.79-2.20) 96.0 (26.79-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.258 , 0.286 0.262 , 0.286	Depositor DCC
R_{free} test set	2103 reflections (3.93%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.0	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	0.92	EDS
Total number of atoms	4104	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.95% 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: NI

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	0.68	6/2054 (0.3%)	1.08	16/2774 (0.6%)
1	B	0.67	8/2043 (0.4%)	1.06	15/2762 (0.5%)
All	All	0.67	14/4097 (0.3%)	1.07	31/5536 (0.6%)

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	1
All	All	0	2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	MSE	SE-CE	-12.29	1.58	1.95
1	A	92	GLU	CB-CG	-9.79	1.23	1.52
1	A	93	MSE	SE-CE	-8.97	1.68	1.95
1	A	91	GLY	CA-C	-8.44	1.42	1.51
1	B	152	MSE	SE-CE	-6.76	1.75	1.95
1	A	92	GLU	CA-CB	6.76	1.64	1.53
1	B	92	GLU	CA-CB	6.17	1.63	1.53
1	B	92	GLU	CB-CG	-6.01	1.34	1.52
1	B	33	MSE	SE-CE	-5.97	1.77	1.95
1	B	196	MSE	SE-CE	-5.96	1.77	1.95
1	B	91	GLY	C-N	5.72	1.41	1.33
1	A	33	MSE	SE-CE	-5.18	1.79	1.95
1	B	91	GLY	N-CA	-5.04	1.40	1.45
1	A	91	GLY	N-CA	-5.03	1.40	1.45

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	GLU	CA-CB-CG	-13.36	87.38	114.10
1	A	91	GLY	CA-C-N	11.50	143.51	121.54
1	A	91	GLY	C-N-CA	11.50	143.51	121.54
1	A	92	GLU	N-CA-CB	10.28	127.86	110.49
1	B	92	GLU	CA-CB-CG	-10.14	93.82	114.10
1	B	91	GLY	CA-C-N	9.60	139.88	121.54
1	B	91	GLY	C-N-CA	9.60	139.88	121.54
1	B	203	THR	N-CA-C	-9.36	94.71	109.50
1	A	91	GLY	O-C-N	9.34	134.25	123.51
1	B	92	GLU	N-CA-CB	8.88	125.50	110.49
1	A	92	GLU	CB-CG-CD	-8.77	97.69	112.60
1	A	91	GLY	N-CA-C	7.63	125.41	110.83
1	A	92	GLU	N-CA-C	-7.61	94.59	110.80
1	B	208	LEU	N-CA-C	7.32	121.49	111.39
1	A	127	ILE	N-CA-C	7.29	118.78	108.36
1	A	16	ASP	N-CA-C	7.23	119.24	111.36
1	B	91	GLY	N-CA-C	6.99	123.94	110.66
1	A	208	LEU	N-CA-C	6.95	120.83	111.24
1	B	91	GLY	O-C-N	6.62	131.12	123.44
1	B	17	ASP	N-CA-C	6.34	122.09	113.97
1	B	127	ILE	N-CA-C	6.29	117.66	108.53
1	B	92	GLU	N-CA-C	-6.29	97.41	110.80
1	B	204	ASP	N-CA-C	-5.92	104.19	112.25
1	B	194	VAL	N-CA-C	5.60	117.69	109.63
1	B	193	HIS	N-CA-C	5.56	117.20	111.03
1	B	152	MSE	N-CA-C	-5.32	98.61	108.24
1	A	169	LEU	N-CA-C	-5.20	102.06	110.32
1	A	92	GLU	CA-C-O	5.17	127.91	120.51
1	A	204	ASP	N-CA-C	-5.15	105.65	112.24
1	A	152	MSE	N-CA-C	-5.10	98.39	107.98
1	A	207	TYR	N-CA-C	5.07	117.39	110.55

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	91	GLY	Mainchain
1	B	91	GLY	Mainchain

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	2015	0	1977	44	0
1	B	2004	0	1953	67	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	41	0	0	3	0
3	B	40	0	0	2	0
All	All	4104	0	3930	110	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 14.

All (110) 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:24:ARG:HG2	1:B:24:ARG:HH11	1.32	0.92
1:B:47:MSE:HE2	1:B:81:LEU:HD12	1.55	0.89
1:A:229:GLN:HE21	1:A:233:LEU:HD13	1.44	0.80
1:B:5:THR:HG22	3:B:370:HOH:O	1.85	0.77
1:B:186:GLN:H	1:B:186:GLN:CD	1.93	0.77
1:A:239:GLU:CD	1:A:239:GLU:H	1.91	0.76
1:B:38:PHE:HB2	1:B:63:HIS:HB2	1.69	0.75
1:B:186:GLN:HB2	1:B:187:PRO:HD3	1.70	0.74
1:B:244:GLN:HE22	1:B:247:LYS:HD3	1.56	0.71
1:A:21:VAL:HG13	1:A:219:GLU:HG3	1.71	0.70
1:B:86:LYS:HE2	3:B:337:HOH:O	1.90	0.70
1:B:24:ARG:HG2	1:B:24:ARG:NH1	2.04	0.70
1:B:7:VAL:HG21	1:B:220:PRO:HG3	1.75	0.69
1:A:219:GLU:HG2	3:A:302:HOH:O	1.91	0.69
1:A:217:ARG:HG3	1:A:217:ARG:HH11	1.58	0.68
1:A:162:ALA:O	1:A:165:VAL:HG22	1.95	0.66
1:B:79:GLU:HG2	1:B:117:LEU:CD1	2.25	0.66
1:B:177:GLY:HA2	1:B:210:PRO:HB3	1.78	0.66
1:B:92:GLU:CG	1:B:128:HIS:HB2	2.26	0.66
1:B:127:ILE:HD12	1:B:140:LEU:HD11	1.77	0.65
1:A:92:GLU:OE2	1:A:128:HIS:ND1	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:PHE:O	1:B:188:LYS:HE2	1.99	0.63
1:B:70:PHE:CD2	1:B:110:VAL:HG22	2.34	0.63
1:B:244:GLN:NE2	1:B:247:LYS:HD3	2.14	0.62
1:A:191:ALA:HB3	1:A:233:LEU:HD23	1.81	0.61
1:A:98:HIS:CB	1:A:130:ARG:HD3	2.31	0.61
1:A:98:HIS:HB2	1:A:130:ARG:HD3	1.84	0.60
1:B:57:TYR:CD1	1:B:86:LYS:HG3	2.37	0.60
1:A:92:GLU:CG	1:A:128:HIS:HB2	2.33	0.59
1:B:79:GLU:HG2	1:B:117:LEU:HD11	1.84	0.59
1:B:229:GLN:HG3	1:B:233:LEU:HD22	1.85	0.59
1:A:253:PHE:O	1:A:254:ASN:CB	2.51	0.57
1:B:7:VAL:HG23	1:B:205:ALA:HB3	1.86	0.57
1:B:244:GLN:HE21	1:B:248:ASN:HD21	1.52	0.57
1:A:40:LYS:O	1:A:44:GLU:HG3	2.04	0.57
1:A:229:GLN:NE2	1:A:233:LEU:HD13	2.18	0.57
1:B:217:ARG:HD2	1:B:218:ASN:O	2.05	0.56
1:A:177:GLY:HA2	1:A:210:PRO:HB3	1.86	0.56
1:B:6:HIS:O	1:B:203:THR:O	2.23	0.56
1:A:141:LEU:HD23	1:A:169:LEU:HD21	1.88	0.56
1:A:161:ILE:O	1:A:165:VAL:HG13	2.07	0.55
1:B:196:MSE:HG3	1:B:244:GLN:HG2	1.89	0.54
1:B:163:ASP:HA	1:B:166:THR:HG22	1.90	0.53
1:A:158:SER:HB2	1:A:159:PRO:HD2	1.90	0.53
1:B:127:ILE:HD11	1:B:152:MSE:CE	2.38	0.53
1:B:92:GLU:HG2	1:B:128:HIS:HB2	1.91	0.53
1:B:75:LEU:C	1:B:75:LEU:HD23	2.34	0.52
1:A:79:GLU:HG3	1:A:121:LEU:HD11	1.90	0.52
1:A:186:GLN:O	1:A:190:VAL:HG23	2.11	0.51
1:B:68:ILE:HD13	1:B:68:ILE:O	2.11	0.51
1:B:79:GLU:HG2	1:B:117:LEU:HD13	1.92	0.51
1:B:253:PHE:O	1:B:254:ASN:CB	2.58	0.50
1:A:186:GLN:HB3	1:A:187:PRO:HD3	1.94	0.50
1:A:100:ASP:OD2	1:B:211:HIS:HE1	1.95	0.49
1:A:160:GLU:O	1:A:164:ILE:HG13	2.12	0.49
1:B:189:GLU:HA	1:B:192:LYS:HZ2	1.77	0.49
1:A:117:LEU:HD22	1:A:121:LEU:HD13	1.93	0.49
1:A:205:ALA:HA	1:A:218:ASN:HB3	1.94	0.49
1:A:38:PHE:HB2	1:A:63:HIS:HB2	1.95	0.48
1:B:11:ASP:OD2	1:B:12:GLU:N	2.46	0.48
1:A:217:ARG:HH11	1:A:217:ARG:CG	2.23	0.48
1:B:176:GLY:O	1:B:179:VAL:HG12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ARG:NH1	1:B:24:ARG:CG	2.73	0.48
1:B:67:ALA:O	1:B:70:PHE:CD1	2.67	0.48
1:B:231:ALA:HB2	1:B:241:VAL:HG21	1.95	0.48
1:A:152:MSE:HE3	3:A:375:HOH:O	2.14	0.47
1:B:246:THR:O	1:B:250:GLU:HG3	2.15	0.47
1:B:194:VAL:CG1	1:B:199:LEU:HD13	2.45	0.46
1:A:55:PHE:CE2	1:A:56:LEU:HD13	2.51	0.46
1:B:77:TRP:NE1	1:B:81:LEU:HD21	2.31	0.45
1:B:203:THR:HB	1:B:218:ASN:ND2	2.31	0.45
1:A:120:ARG:O	1:A:122:LYS:HD2	2.15	0.45
1:B:126:ILE:HD12	1:B:126:ILE:N	2.31	0.45
1:A:239:GLU:CD	1:A:239:GLU:N	2.68	0.45
1:B:127:ILE:HG13	1:B:152:MSE:SE	2.67	0.45
1:B:229:GLN:HG3	1:B:233:LEU:CD2	2.46	0.45
1:A:217:ARG:HG3	1:A:217:ARG:O	2.15	0.45
1:B:38:PHE:CD2	1:B:63:HIS:HB2	2.52	0.45
1:B:203:THR:HB	1:B:218:ASN:HD21	1.82	0.45
1:B:39:ASN:HD22	1:B:39:ASN:HA	1.52	0.45
1:A:129:ASN:HB3	3:A:383:HOH:O	2.16	0.45
1:B:127:ILE:HD11	1:B:152:MSE:HE1	1.98	0.45
1:B:127:ILE:HD11	1:B:152:MSE:SE	2.67	0.44
1:B:120:ARG:O	1:B:122:LYS:HE2	2.16	0.44
1:B:205:ALA:HA	1:B:218:ASN:HB3	1.98	0.44
1:A:178:PRO:HA	1:A:181:PHE:CD2	2.53	0.44
1:B:12:GLU:C	1:B:14:TYR:H	2.26	0.44
1:A:84:HIS:CE1	1:A:86:LYS:HB2	2.53	0.44
1:A:39:ASN:HD22	1:A:39:ASN:HA	1.54	0.43
1:B:64:PRO:HB3	1:B:94:GLY:O	2.18	0.43
1:B:239:GLU:CD	1:B:239:GLU:H	2.25	0.43
1:A:217:ARG:CG	1:A:217:ARG:NH1	2.81	0.43
1:A:48:LYS:HB3	1:A:48:LYS:HE2	1.84	0.42
1:A:117:LEU:C	1:A:117:LEU:HD13	2.44	0.42
1:B:70:PHE:CG	1:B:110:VAL:HG22	2.54	0.42
1:B:67:ALA:O	1:B:70:PHE:HD1	2.03	0.42
1:B:153:HIS:HA	1:B:174:SER:HB3	2.02	0.42
1:B:144:HIS:HA	1:B:146:GLU:OE1	2.19	0.42
1:A:63:HIS:ND1	1:A:64:PRO:HD2	2.35	0.41
1:B:53:TYR:HB2	1:B:56:LEU:HD22	2.01	0.41
1:B:2:LEU:HD11	1:B:201:VAL:HG21	2.02	0.41
1:B:14:TYR:OH	1:B:205:ALA:HB1	2.21	0.41
1:B:186:GLN:CD	1:B:186:GLN:N	2.71	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:VAL:HG11	1:B:199:LEU:HD13	2.01	0.41
1:A:110:VAL:HG12	1:A:114:GLN:HE21	1.86	0.41
1:A:205:ALA:CB	1:A:218:ASN:HB3	2.51	0.41
1:B:93:MSE:CG	1:B:127:ILE:HG22	2.51	0.41
1:B:70:PHE:CG	1:B:70:PHE:O	2.74	0.40
1:A:7:VAL:O	1:A:35:VAL:HA	2.21	0.40
1:A:172:TYR:CZ	1:A:198:ARG:NH2	2.89	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	251/265 (95%)	243 (97%)	8 (3%)	0	100	100
1	B	251/265 (95%)	241 (96%)	10 (4%)	0	100	100
All	All	502/530 (95%)	484 (96%)	18 (4%)	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	216/227 (95%)	207 (96%)	9 (4%)	26	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	213/227 (94%)	197 (92%)	16 (8%)	12	14
All	All	429/454 (94%)	404 (94%)	25 (6%)	18	22

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	56	LEU
1	A	153	HIS
1	A	165	VAL
1	A	179	VAL
1	A	211	HIS
1	A	217	ARG
1	A	233	LEU
1	A	239	GLU
1	B	5	THR
1	B	17	ASP
1	B	24	ARG
1	B	39	ASN
1	B	51	ASP
1	B	56	LEU
1	B	68	ILE
1	B	86	LYS
1	B	95	LEU
1	B	127	ILE
1	B	153	HIS
1	B	166	THR
1	B	175	LEU
1	B	199	LEU
1	B	233	LEU
1	B	239	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	74	HIS
1	A	134	GLN
1	A	167	ASN
1	A	183	ASN
1	A	229	GLN

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Mol	Chain	Res	Type
1	A	244	GLN
1	A	248	ASN
1	B	39	ASN
1	B	74	HIS
1	B	134	GLN
1	B	167	ASN
1	B	183	ASN
1	B	211	HIS
1	B	244	GLN
1	B	248	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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 [i](#)

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 [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	248/265 (93%)	0.36	7 (2%) 55 52	12, 28, 39, 47	0
1	B	248/265 (93%)	0.47	12 (4%) 35 32	16, 30, 42, 49	0
All	All	496/530 (93%)	0.42	19 (3%) 44 41	12, 29, 42, 49	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	70	PHE	5.1
1	A	20	GLU	4.6
1	B	254	ASN	4.5
1	A	254	ASN	4.4
1	B	204	ASP	4.3
1	B	17	ASP	3.3
1	B	207	TYR	3.0
1	A	105	ASP	3.0
1	A	207	TYR	2.8
1	A	168	LYS	2.7
1	B	39	ASN	2.6
1	B	16	ASP	2.5
1	B	186	GLN	2.4
1	A	99	TRP	2.3
1	B	160	GLU	2.2
1	B	13	GLN	2.1
1	B	91	GLY	2.1
1	B	158	SER	2.1
1	A	136	CYS	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	NI	B	298	1/1	0.12	0.35	30,30,30,30	0
2	NI	A	299	1/1	0.41	0.25	22,22,22,22	0
2	NI	A	300	1/1	0.97	0.04	15,15,15,15	0
2	NI	B	297	1/1	0.98	0.03	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.