



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:14 PM UTC

PDB ID : 1YRT / pdb_00001yrt
Title : Crystal Structure analysis of the adenylyl cyclase catalytic domain of adenylyl cyclase toxin of Bordetella pertussis in presence of c-terminal calmodulin
Authors : Guo, Q.; Shen, Y.; Lee, Y.S.; Gibbs, C.S.; Mrksich, M.; Tang, W.J.
Deposited on : 2005-02-04
Resolution : 2.10 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

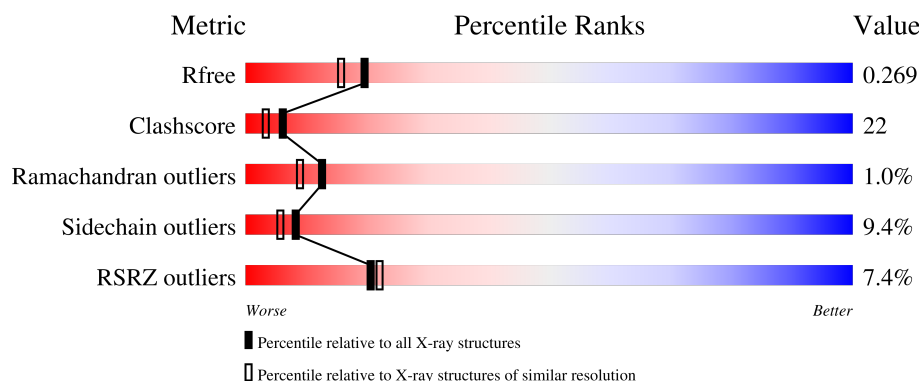
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

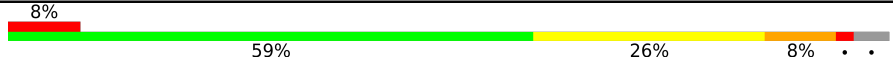
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	
2	B	74	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3397 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 Bifunctional hemolysin-adenylate cyclase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2684	1664	495	519	6			

- Molecule 2 is a protein called Calmodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	69	Total	C	N	O	S	0	0	0
			551	335	90	122	4			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Ca	0	0
			2	2		

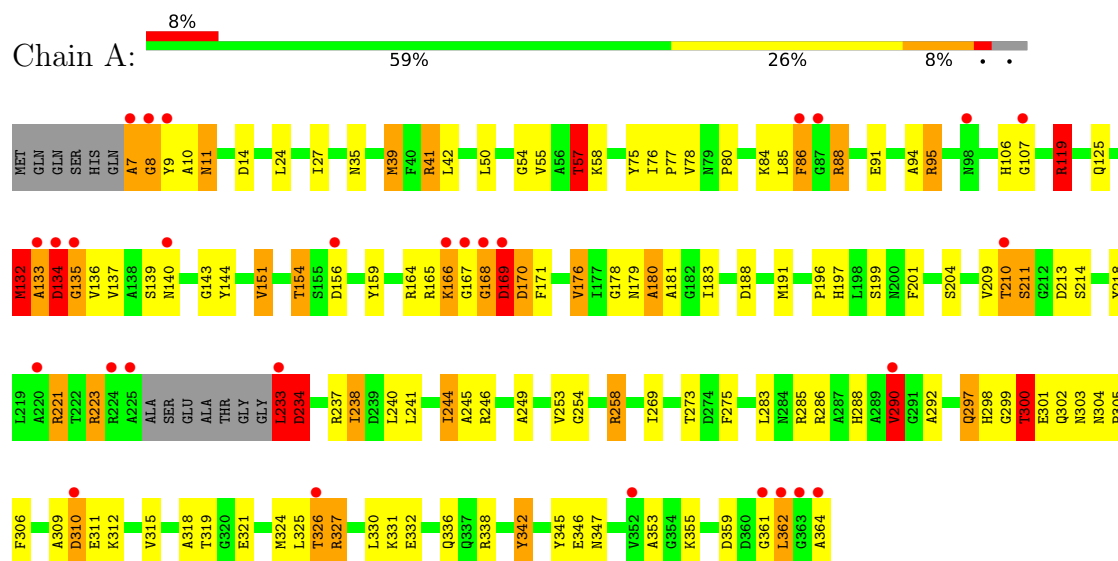
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	133	Total	O	0	0
			133	133		
4	B	27	Total	O	0	0
			27	27		

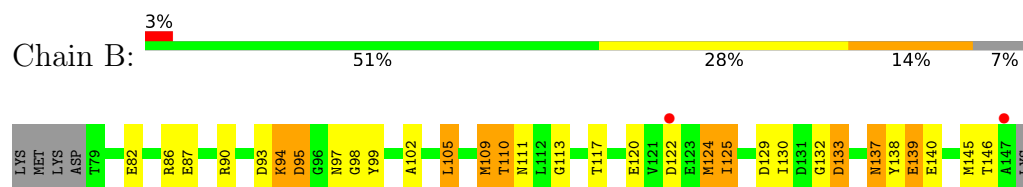
3 Residue-property plots [i](#)

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: Bifunctional hemolysin-adenylate cyclase



- Molecule 2: Calmodulin



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	79.36Å 79.36Å 139.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.10) 96.5 (30.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.10Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.220 , 0.270 0.221 , 0.269	Depositor DCC
R_{free} test set	1277 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3397	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 5.72% 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: CA

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	1.54	13/2728 (0.5%)	1.78	67/3682 (1.8%)
2	B	1.39	3/556 (0.5%)	1.84	18/746 (2.4%)
All	All	1.52	16/3284 (0.5%)	1.79	85/4428 (1.9%)

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	9

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	132	MET	C-N	-34.05	0.89	1.33
1	A	135	GLY	N-CA	26.94	1.84	1.45
1	A	8	GLY	C-N	15.51	1.55	1.33
1	A	7	ALA	C-N	-14.39	1.12	1.33
1	A	9	TYR	N-CA	9.19	1.58	1.46
2	B	124	MET	SD-CE	-7.73	1.60	1.79
1	A	10	ALA	N-CA	6.52	1.55	1.46
1	A	191	MET	SD-CE	6.30	1.95	1.79
1	A	234	ASP	C-N	6.19	1.42	1.33
2	B	110	THR	N-CA	5.99	1.54	1.46
1	A	168	GLY	C-N	-5.88	1.25	1.33
1	A	167	GLY	CA-C	5.79	1.59	1.51
1	A	39	MET	SD-CE	-5.56	1.65	1.79
2	B	110	THR	CA-CB	5.21	1.62	1.53
1	A	57	THR	CB-CG2	-5.03	1.35	1.52
1	A	249	ALA	CA-CB	5.00	1.61	1.53

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	ALA	O-C-N	-32.68	70.72	123.00
1	A	10	ALA	N-CA-C	23.18	140.74	113.15
2	B	110	THR	N-CA-CB	-21.59	76.75	110.44
1	A	327	ARG	N-CA-CB	-19.33	81.73	110.53
1	A	234	ASP	CA-CB-CG	-17.12	95.48	112.60
1	A	10	ALA	N-CA-CB	-14.88	87.55	110.46
1	A	326	THR	N-CA-C	-14.35	87.40	109.79
1	A	233	LEU	CA-C-N	13.88	148.04	121.54
1	A	233	LEU	C-N-CA	13.88	148.04	121.54
2	B	95	ASP	N-CA-C	-13.68	92.89	110.53
1	A	210	THR	N-CA-C	-11.76	99.14	108.78
1	A	311	GLU	N-CA-C	11.22	126.34	112.87
1	A	134	ASP	CA-C-N	-10.90	100.05	121.41
1	A	134	ASP	C-N-CA	-10.90	100.05	121.41
2	B	133	ASP	N-CA-C	-9.99	96.57	110.35
2	B	133	ASP	CA-C-N	-9.16	103.45	121.41
2	B	133	ASP	C-N-CA	-9.16	103.45	121.41
1	A	180	ALA	N-CA-C	-9.13	101.40	112.54
1	A	290	VAL	CB-CA-C	8.94	125.95	111.29
1	A	290	VAL	CA-C-N	-8.67	104.42	121.41
1	A	290	VAL	C-N-CA	-8.67	104.42	121.41
1	A	211	SER	N-CA-C	-8.63	97.72	110.06
1	A	290	VAL	O-C-N	-8.53	111.90	122.57
1	A	210	THR	CA-C-O	8.38	122.80	117.94
1	A	132	MET	O-C-N	-8.10	113.41	122.72
1	A	233	LEU	O-C-N	-8.09	110.05	123.00
2	B	110	THR	N-CA-C	-7.33	104.14	113.23
2	B	111	ASN	N-CA-C	-7.20	103.43	111.28
1	A	86	PHE	CA-C-N	-7.03	107.64	121.41
1	A	86	PHE	C-N-CA	-7.03	107.64	121.41
2	B	95	ASP	CA-C-N	-6.95	107.78	121.41
2	B	95	ASP	C-N-CA	-6.95	107.78	121.41
1	A	7	ALA	CA-C-N	6.91	134.95	121.41
1	A	7	ALA	C-N-CA	6.91	134.95	121.41
1	A	309	ALA	N-CA-C	-6.84	102.23	110.44
1	A	119	ARG	N-CA-C	-6.84	103.89	111.82
1	A	210	THR	CB-CA-C	-6.71	107.40	117.07
1	A	345	TYR	N-CA-C	-6.65	99.00	109.50
1	A	290	VAL	N-CA-CB	-6.60	100.34	111.23
1	A	132	MET	CA-C-N	6.54	132.81	122.60
1	A	132	MET	C-N-CA	6.54	132.81	122.60
1	A	119	ARG	CB-CG-CD	6.51	126.26	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ARG	N-CA-C	6.41	119.21	110.55
1	A	300	THR	N-CA-CB	-6.33	100.83	110.45
1	A	57	THR	N-CA-CB	-6.33	99.82	110.83
1	A	132	MET	N-CA-C	6.24	116.47	108.24
1	A	327	ARG	N-CA-C	-6.19	105.77	113.38
1	A	86	PHE	CA-CB-CG	-6.17	107.63	113.80
1	A	77	PRO	N-CA-C	6.11	120.78	111.19
1	A	290	VAL	N-CA-C	-6.10	96.66	109.34
1	A	86	PHE	CA-C-O	-6.05	113.94	121.02
1	A	169	ASP	CA-C-N	-6.04	112.17	122.33
1	A	169	ASP	C-N-CA	-6.04	112.17	122.33
1	A	41	ARG	N-CA-C	-6.03	101.56	110.48
2	B	94	LYS	N-CA-C	5.96	118.26	111.11
1	A	86	PHE	O-C-N	-5.91	115.05	122.20
2	B	109	MET	N-CA-C	5.90	117.71	111.28
1	A	234	ASP	N-CA-CB	5.82	120.32	110.49
1	A	214	SER	N-CA-C	-5.76	102.88	110.43
1	A	119	ARG	NE-CZ-NH1	5.74	127.23	121.50
1	A	310	ASP	CA-C-N	5.64	131.49	120.99
1	A	310	ASP	C-N-CA	5.64	131.49	120.99
1	A	292	ALA	N-CA-C	5.61	118.82	110.52
1	A	300	THR	OG1-CB-CG2	5.61	120.51	109.30
1	A	9	TYR	CA-C-N	-5.57	111.18	122.31
1	A	9	TYR	C-N-CA	-5.57	111.18	122.31
1	A	258	ARG	NE-CZ-NH2	-5.54	114.21	119.20
2	B	109	MET	O-C-N	-5.50	116.30	122.12
1	A	154	THR	N-CA-C	-5.42	101.98	110.17
1	A	55	VAL	N-CA-C	-5.42	102.59	109.58
1	A	54	GLY	N-CA-C	5.39	123.86	115.72
2	B	113	GLY	N-CA-C	5.36	123.16	114.90
2	B	130	ILE	N-CA-C	5.29	115.92	110.36
2	B	130	ILE	CB-CA-C	-5.27	105.14	111.88
1	A	258	ARG	NE-CZ-NH1	5.22	126.72	121.50
1	A	169	ASP	N-CA-C	5.22	121.91	110.80
1	A	234	ASP	O-C-N	-5.14	115.76	122.59
1	A	209	VAL	CA-C-N	-5.12	117.51	123.04
1	A	209	VAL	C-N-CA	-5.12	117.51	123.04
2	B	133	ASP	O-C-N	-5.10	115.45	122.23
1	A	201	PHE	N-CA-C	5.08	119.44	112.88
1	A	178	GLY	CA-C-N	5.08	129.46	120.87
1	A	178	GLY	C-N-CA	5.08	129.46	120.87
2	B	109	MET	CA-C-N	-5.07	112.75	121.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	109	MET	C-N-CA	-5.07	112.75	121.66

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	MET	Mainchain
1	A	133	ALA	Mainchain
1	A	134	ASP	Peptide
1	A	233	LEU	Peptide
1	A	342	TYR	Sidechain
1	A	7	ALA	Mainchain,Peptide
1	A	8	GLY	Mainchain,Peptide

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	2684	0	2622	121	1
2	B	551	0	503	29	0
3	B	2	0	0	0	0
4	A	133	0	0	2	1
4	B	27	0	0	2	0
All	All	3397	0	3125	143	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 22.

All (143) 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:135:GLY:N	1:A:135:GLY:CA	1.84	1.39
1:A:169:ASP:OD2	1:A:169:ASP:O	1.65	1.14
1:A:133:ALA:HB3	1:A:136:VAL:H	1.05	1.07
1:A:210:THR:HG22	1:A:211:SER:H	0.96	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ALA:HB2	1:A:136:VAL:HB	1.39	1.04
1:A:210:THR:HG22	1:A:211:SER:N	1.72	1.03
1:A:133:ALA:HB3	1:A:136:VAL:N	1.75	1.02
1:A:210:THR:CG2	1:A:211:SER:H	1.70	1.00
1:A:210:THR:HG21	1:A:213:ASP:HB2	1.43	0.99
2:B:86:ARG:HH11	2:B:86:ARG:HG2	1.32	0.92
1:A:134:ASP:C	1:A:135:GLY:CA	2.46	0.87
1:A:223:ARG:NE	1:A:223:ARG:HA	1.90	0.85
1:A:210:THR:HG23	1:A:218:TYR:CD1	2.12	0.85
1:A:133:ALA:CB	1:A:136:VAL:HB	2.07	0.84
1:A:210:THR:HG23	1:A:218:TYR:HD1	1.43	0.84
1:A:223:ARG:HA	1:A:223:ARG:HE	1.45	0.80
1:A:179:ASN:HB2	1:A:183:ILE:H	1.46	0.79
2:B:137:ASN:C	2:B:137:ASN:HD22	1.90	0.79
1:A:210:THR:HG21	1:A:213:ASP:CB	2.13	0.79
1:A:132:MET:HE2	1:A:135:GLY:N	1.98	0.78
1:A:297:GLN:HE21	1:A:297:GLN:HA	1.48	0.78
1:A:134:ASP:O	1:A:134:ASP:OD1	2.02	0.77
2:B:109:MET:HG3	2:B:124:MET:HE1	1.68	0.74
1:A:302:GLN:NE2	1:A:347:ASN:H	1.87	0.72
2:B:93:ASP:OD1	2:B:97:ASN:O	2.07	0.72
1:A:223:ARG:HE	1:A:223:ARG:CA	2.01	0.71
1:A:91:GLU:OE2	1:A:95:ARG:HG3	1.89	0.71
1:A:300:THR:HG23	1:A:302:GLN:H	1.56	0.71
1:A:362:LEU:HG	2:B:90:ARG:NH2	2.05	0.71
1:A:237:ARG:NH1	1:A:241:LEU:HD11	2.06	0.70
2:B:105:LEU:HD13	2:B:124:MET:HE3	1.73	0.70
1:A:41:ARG:HE	1:A:310:ASP:CG	2.00	0.69
1:A:151:VAL:HG22	1:A:159:TYR:HB3	1.73	0.69
1:A:298:HIS:HD2	1:A:299:GLY:O	1.76	0.68
1:A:169:ASP:OD2	1:A:169:ASP:C	2.36	0.68
2:B:86:ARG:HG2	2:B:86:ARG:NH1	2.08	0.68
1:A:133:ALA:CB	1:A:136:VAL:CB	2.71	0.68
1:A:327:ARG:HA	1:A:330:LEU:HB3	1.75	0.67
1:A:362:LEU:HG	2:B:90:ARG:HH22	1.60	0.67
1:A:300:THR:CG2	1:A:302:GLN:H	2.07	0.67
1:A:286:ARG:HD3	4:A:368:HOH:O	1.95	0.67
1:A:41:ARG:HH21	1:A:310:ASP:HB3	1.62	0.64
1:A:57:THR:HG23	1:A:188:ASP:HB3	1.80	0.63
1:A:237:ARG:HH12	1:A:241:LEU:HD11	1.65	0.62
1:A:325:LEU:C	1:A:326:THR:O	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:ARG:HD2	4:B:807:HOH:O	1.99	0.62
1:A:154:THR:OG1	1:A:156:ASP:HB3	1.99	0.62
2:B:137:ASN:ND2	2:B:140:GLU:H	1.98	0.62
2:B:86:ARG:HH11	2:B:86:ARG:CG	2.04	0.61
1:A:246:ARG:HB2	2:B:145:MET:HE1	1.82	0.61
1:A:300:THR:HG21	4:A:380:HOH:O	1.99	0.61
1:A:179:ASN:HB2	1:A:183:ILE:N	2.15	0.60
1:A:285:ARG:HD2	1:A:285:ARG:C	2.26	0.60
2:B:137:ASN:C	2:B:137:ASN:ND2	2.59	0.60
1:A:11:ASN:ND2	1:A:14:ASP:H	2.00	0.60
2:B:137:ASN:HD21	2:B:139:GLU:HG3	1.67	0.60
1:A:39:MET:HG2	1:A:315:VAL:HG22	1.83	0.59
1:A:253:VAL:O	1:A:253:VAL:HG12	2.01	0.59
1:A:364:ALA:HA	2:B:138:TYR:CE2	2.37	0.59
1:A:85:LEU:O	1:A:88:ARG:HG2	2.04	0.58
1:A:151:VAL:CG2	1:A:159:TYR:HB3	2.32	0.58
1:A:164:ARG:HD3	1:A:171:PHE:CE2	2.39	0.58
1:A:304:ASN:ND2	1:A:306:PHE:H	2.01	0.58
1:A:27:ILE:HG12	1:A:283:LEU:HD22	1.86	0.57
1:A:253:VAL:HG12	1:A:258:ARG:HH21	1.69	0.57
1:A:135:GLY:N	1:A:135:GLY:C	2.62	0.57
1:A:133:ALA:O	1:A:134:ASP:HB3	2.04	0.56
1:A:213:ASP:OD2	1:A:221:ARG:NH1	2.39	0.56
2:B:102:ALA:HA	2:B:125:ILE:CD1	2.36	0.56
1:A:137:VAL:HG23	1:A:144:TYR:HE2	1.69	0.56
1:A:302:GLN:HE21	1:A:347:ASN:H	1.53	0.56
1:A:132:MET:HA	1:A:136:VAL:O	2.06	0.55
1:A:332:GLU:O	1:A:336:GLN:HG3	2.06	0.55
1:A:168:GLY:O	1:A:170:ASP:N	2.40	0.55
1:A:210:THR:CG2	1:A:211:SER:N	2.41	0.55
1:A:364:ALA:HB2	2:B:82:GLU:HG3	1.90	0.54
2:B:129:ASP:OD1	2:B:133:ASP:O	2.25	0.54
2:B:117:THR:OG1	2:B:120:GLU:HG3	2.08	0.54
2:B:137:ASN:HD21	2:B:140:GLU:H	1.56	0.53
2:B:97:ASN:O	2:B:99:TYR:N	2.40	0.53
1:A:244:ILE:HG13	1:A:245:ALA:N	2.22	0.53
1:A:210:THR:CG2	1:A:213:ASP:H	2.22	0.53
1:A:223:ARG:HE	1:A:223:ARG:C	2.16	0.53
1:A:297:GLN:HE21	1:A:297:GLN:CA	2.21	0.52
1:A:50:LEU:HD21	1:A:119:ARG:HD2	1.91	0.52
1:A:269:ILE:CG2	1:A:298:HIS:HA	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:361:GLY:O	1:A:362:LEU:HB2	2.11	0.51
1:A:197:HIS:HD2	1:A:199:SER:OG	1.94	0.51
1:A:304:ASN:HD22	1:A:305:PRO:HD2	1.75	0.51
2:B:95:ASP:OD2	2:B:97:ASN:ND2	2.44	0.51
1:A:218:TYR:CD2	1:A:240:LEU:HD21	2.46	0.50
2:B:86:ARG:NH1	2:B:86:ARG:CG	2.66	0.50
1:A:301:GLU:O	1:A:301:GLU:HG3	2.12	0.50
1:A:210:THR:HG21	1:A:213:ASP:H	1.76	0.49
1:A:42:LEU:CD2	1:A:324:MET:HE1	2.43	0.49
1:A:164:ARG:HH11	1:A:171:PHE:HE2	1.59	0.49
1:A:78:VAL:O	1:A:80:PRO:HD3	2.12	0.49
1:A:285:ARG:HD2	1:A:285:ARG:O	2.13	0.48
1:A:75:TYR:HB3	1:A:176:VAL:HG21	1.95	0.48
1:A:258:ARG:HH22	2:B:87:GLU:CD	2.22	0.47
1:A:258:ARG:HD2	1:A:303:ASN:OD1	2.14	0.47
1:A:57:THR:HG23	1:A:188:ASP:CB	2.44	0.47
1:A:58:LYS:HD3	1:A:76:ILE:HD11	1.97	0.46
1:A:269:ILE:HG23	1:A:298:HIS:HA	1.97	0.46
1:A:302:GLN:HE21	1:A:346:GLU:HA	1.79	0.46
1:A:107:GLY:O	1:A:180:ALA:HA	2.15	0.46
1:A:86:PHE:CZ	1:A:143:GLY:HA2	2.51	0.46
2:B:93:ASP:CG	2:B:97:ASN:O	2.59	0.46
1:A:134:ASP:O	1:A:135:GLY:CA	2.65	0.45
1:A:197:HIS:CD2	1:A:199:SER:H	2.33	0.45
1:A:254:GLY:HA3	1:A:273:THR:HG22	1.97	0.45
1:A:253:VAL:HG11	2:B:87:GLU:HB3	1.98	0.45
1:A:269:ILE:HG22	1:A:269:ILE:O	2.15	0.45
1:A:253:VAL:O	1:A:253:VAL:CG1	2.64	0.45
1:A:11:ASN:HD22	1:A:14:ASP:H	1.66	0.44
1:A:234:ASP:O	1:A:238:ILE:HD12	2.17	0.44
1:A:179:ASN:HB3	1:A:181:ALA:H	1.82	0.44
1:A:288:HIS:O	1:A:290:VAL:O	2.36	0.44
1:A:304:ASN:HD22	1:A:305:PRO:CD	2.31	0.44
1:A:137:VAL:HG23	1:A:144:TYR:CE2	2.52	0.43
1:A:57:THR:HG22	1:A:297:GLN:CG	2.49	0.43
1:A:35:ASN:HD22	1:A:318:ALA:HB1	1.83	0.43
1:A:319:THR:OG1	1:A:321:GLU:HG3	2.19	0.43
1:A:139:SER:HB3	1:A:144:TYR:CG	2.54	0.43
1:A:165:ARG:O	1:A:166:LYS:C	2.62	0.42
1:A:106:HIS:ND1	1:A:106:HIS:N	2.67	0.42
2:B:146:THR:O	2:B:146:THR:HG22	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:ALA:O	1:A:355:LYS:HG2	2.20	0.42
1:A:35:ASN:ND2	1:A:342:TYR:OH	2.48	0.42
1:A:346:GLU:O	1:A:347:ASN:C	2.63	0.42
1:A:133:ALA:HB3	1:A:136:VAL:CA	2.46	0.41
2:B:132:GLY:C	2:B:133:ASP:O	2.61	0.41
1:A:41:ARG:HH21	1:A:310:ASP:CB	2.32	0.41
1:A:75:TYR:HB3	1:A:176:VAL:CG2	2.50	0.41
2:B:95:ASP:OD1	2:B:95:ASP:C	2.63	0.41
2:B:95:ASP:HB3	4:B:820:HOH:O	2.19	0.41
1:A:197:HIS:HD2	1:A:199:SER:H	1.67	0.41
1:A:196:PRO:HB3	1:A:275:PHE:CE2	2.56	0.41
1:A:302:GLN:HE21	1:A:347:ASN:N	2.17	0.41
1:A:331:LYS:HB3	1:A:331:LYS:HE2	1.79	0.41
1:A:133:ALA:CB	1:A:136:VAL:HG23	2.50	0.40
1:A:91:GLU:O	1:A:94:ALA:HB3	2.22	0.40
1:A:269:ILE:HG23	1:A:269:ILE:HD12	1.77	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)
1:A:233:LEU:CD2	4:A:431:HOH:O[6_555]	1.96	0.24

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	347/364 (95%)	327 (94%)	17 (5%)	3 (1%)	14	10
2	B	67/74 (90%)	64 (96%)	2 (3%)	1 (2%)	8	4
All	All	414/438 (94%)	391 (94%)	19 (5%)	4 (1%)	12	9

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	ASP
1	A	234	ASP
1	A	134	ASP
2	B	98	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	271/280 (97%)	247 (91%)	24 (9%)	9	6
2	B	59/64 (92%)	52 (88%)	7 (12%)	5	3
All	All	330/344 (96%)	299 (91%)	31 (9%)	8	6

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	24	LEU
1	A	57	THR
1	A	84	LYS
1	A	95	ARG
1	A	119	ARG
1	A	125	GLN
1	A	140	ASN
1	A	151	VAL
1	A	166	LYS
1	A	169	ASP
1	A	170	ASP
1	A	176	VAL
1	A	204	SER
1	A	221	ARG
1	A	223	ARG
1	A	238	ILE
1	A	244	ILE
1	A	290	VAL

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Mol	Chain	Res	Type
1	A	297	GLN
1	A	300	THR
1	A	312	LYS
1	A	359	ASP
1	A	362	LEU
2	B	94	LYS
2	B	105	LEU
2	B	110	THR
2	B	122	ASP
2	B	125	ILE
2	B	137	ASN
2	B	139	GLU

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	11	ASN
1	A	35	ASN
1	A	108	HIS
1	A	197	HIS
1	A	260	GLN
1	A	297	GLN
1	A	298	HIS
1	A	302	GLN
1	A	304	ASN
1	A	337	GLN
2	B	111	ASN
2	B	137	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 [i](#)

Of 2 ligands modelled in this entry, 2 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 [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	7:ALA	C	8:GLY	N	1.12
1	A	132:MET	C	133:ALA	N	0.89

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	351/364 (96%)	0.34	29 (8%) 17 18	13, 25, 46, 56	0
2	B	69/74 (93%)	0.16	2 (2%) 53 56	14, 23, 38, 49	0
All	All	420/438 (95%)	0.31	31 (7%) 20 22	13, 25, 46, 56	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	GLY	5.4
1	A	7	ALA	5.0
1	A	233	LEU	5.0
1	A	135	GLY	4.1
1	A	364	ALA	3.9
1	A	362	LEU	3.9
1	A	363	GLY	3.6
1	A	168	GLY	3.6
1	A	156	ASP	3.6
1	A	310	ASP	3.6
1	A	134	ASP	3.4
1	A	224	ARG	2.9
2	B	147	ALA	2.9
2	B	122	ASP	2.9
1	A	290	VAL	2.8
1	A	167	GLY	2.7
1	A	210	THR	2.7
1	A	352	VAL	2.6
1	A	133	ALA	2.5
1	A	361	GLY	2.5
1	A	169	ASP	2.4
1	A	107	GLY	2.4
1	A	225	ALA	2.3
1	A	86	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	140	ASN	2.2
1	A	166	LYS	2.2
1	A	9	TYR	2.2
1	A	98	ASN	2.1
1	A	87	GLY	2.1
1	A	220	ALA	2.1
1	A	326	THR	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
3	CA	B	800	1/1	0.99	0.02	21,21,21,21	0
3	CA	B	801	1/1	0.99	0.03	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.