Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 11:52 PM GMT

PDB ID	:	1YRT
Title	:	Crystal Structure analysis of the adenylyl cyclaes 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

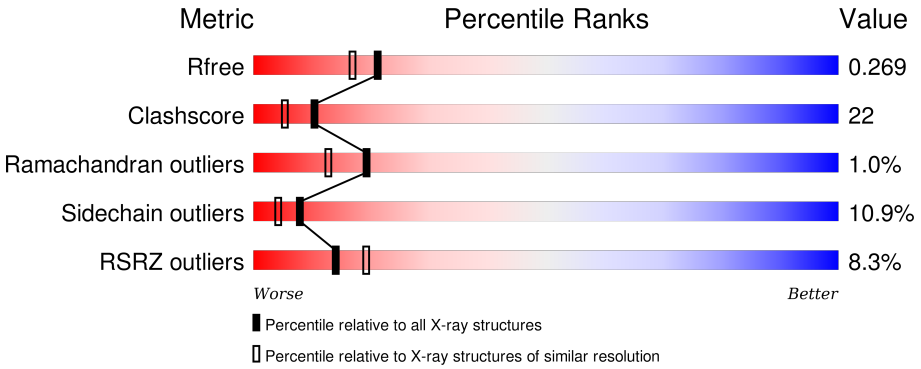
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (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 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 <=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	9% 63% 24% 6%
2	B	74	3% 59% 24% 8% 7%

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 (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is water.

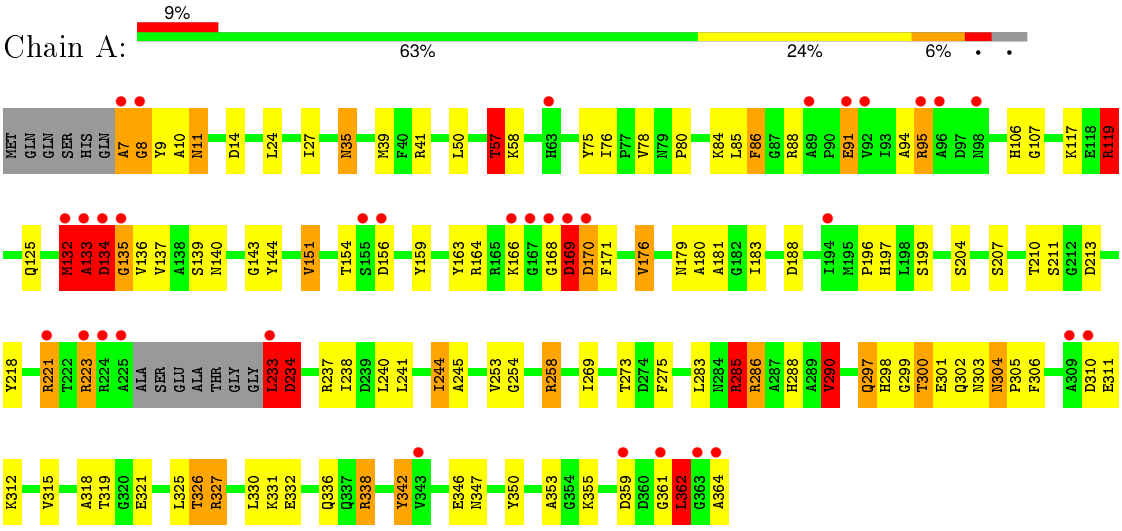
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	133	Total	O	0	0
			133	133		
4	B	27	Total	O	0	0
			27	27		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA 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 (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.

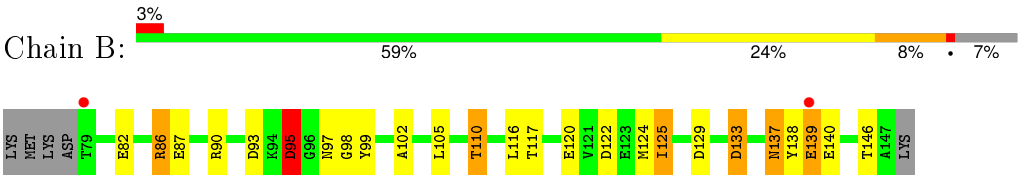
- Molecule 1: Bifunctional hemolysin-adenylate cyclase

Chain A:



- Molecule 2: Calmodulin

Chain B:



4 Data and refinement statistics (i)

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 26.16 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.10) 96.6 (26.16-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [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	1244 reflections (4.83%)	DCC
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 , 50.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 26352 reflections	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	1.22	7/2728 (0.3%)	1.45	35/3682 (1.0%)
2	B	1.09	0/556	1.35	7/746 (0.9%)
All	All	1.20	7/3284 (0.2%)	1.43	42/4428 (0.9%)

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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	135	GLY	N-CA	25.59	1.84	1.46
1	A	132	MET	C-N	-19.25	0.89	1.34
1	A	7	ALA	C-N	-11.43	1.12	1.33
1	A	8	GLY	C-N	9.28	1.55	1.34
1	A	91	GLU	CB-CG	6.34	1.64	1.52
1	A	9	TYR	N-CA	5.84	1.58	1.46
1	A	350	TYR	CD2-CE2	5.70	1.48	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	7	ALA	O-C-N	-30.87	70.72	123.20
1	A	7	ALA	CA-C-N	19.07	154.34	116.20
2	B	110	THR	N-CA-CB	-17.66	76.75	110.30
1	A	10	ALA	N-CA-CB	-16.11	87.55	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	327	ARG	N-CA-CB	-16.04	81.73	110.60
1	A	119	ARG	NE-CZ-NH1	13.87	127.23	120.30
1	A	258	ARG	NE-CZ-NH1	12.84	126.72	120.30
1	A	258	ARG	NE-CZ-NH2	-12.18	114.21	120.30
1	A	10	ALA	N-CA-C	11.01	140.74	111.00
1	A	119	ARG	NE-CZ-NH2	-10.75	114.92	120.30
1	A	134	ASP	C-N-CA	-10.59	100.05	122.30
1	A	233	LEU	C-N-CA	10.54	148.04	121.70
1	A	338	ARG	NE-CZ-NH2	-10.53	115.03	120.30
1	A	338	ARG	NE-CZ-NH1	10.50	125.55	120.30
1	A	8	GLY	CA-C-N	-9.51	96.27	117.20
2	B	133	ASP	C-N-CA	-8.97	103.45	122.30
1	A	326	THR	N-CA-C	-8.74	87.40	111.00
1	A	290	VAL	C-N-CA	-8.52	104.42	122.30
1	A	234	ASP	CA-CB-CG	-8.14	95.48	113.40
1	A	233	LEU	O-C-N	-7.91	110.05	122.70
1	A	290	VAL	CB-CA-C	7.66	125.95	111.40
1	A	86	PHE	CA-C-N	7.35	130.91	116.20
1	A	86	PHE	C-N-CA	-6.98	107.64	122.30
2	B	95	ASP	C-N-CA	-6.91	107.78	122.30
2	B	95	ASP	N-CA-C	-6.71	92.89	111.00
1	A	290	VAL	O-C-N	-6.65	111.90	123.20
1	A	7	ALA	C-N-CA	6.03	134.95	122.30
1	A	290	VAL	CA-C-N	5.93	128.05	116.20
1	A	132	MET	O-C-N	-5.81	113.41	122.70
1	A	311	GLU	N-CA-C	5.68	126.34	111.00
1	A	119	ARG	CB-CG-CD	5.64	126.26	111.60
1	A	57	THR	N-CA-CB	-5.52	99.82	110.30
1	A	285	ARG	NE-CZ-NH1	-5.46	117.57	120.30
2	B	133	ASP	CB-CG-OD1	5.46	123.21	118.30
1	A	286	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	A	234	ASP	N-CA-CB	5.40	120.32	110.60
2	B	133	ASP	N-CA-C	-5.35	96.57	111.00
1	A	290	VAL	N-CA-C	-5.31	96.66	111.00
2	B	86	ARG	NE-CZ-NH1	-5.18	117.71	120.30
1	A	133	ALA	CA-C-N	-5.16	105.84	117.20
1	A	290	VAL	N-CA-CB	-5.07	100.34	111.50
1	A	362	LEU	CA-CB-CG	5.00	126.80	115.30

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 (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	2684	0	2621	117	1
2	B	551	0	503	26	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	3124	137	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 (137) 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:O	1:A:169:ASP:OD2	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
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:223:ARG:HA	1:A:223:ARG:NE	1.90	0.85
1:A:210:THR:HG23	1:A:218:TYR:CD1	2.12	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:134:ASP:C	1:A:135:GLY:CA	2.46	0.82
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
1:A:210:THR:HG21	1:A:213:ASP:CB	2.13	0.79
1:A:297:GLN:HA	1:A:297:GLN:HE21	1.48	0.78
1:A:134:ASP:O	1:A:134:ASP:OD1	2.02	0.77
2:B:137:ASN:C	2:B:137:ASN:HD22	1.90	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:CA	1:A:223:ARG:HE	2.01	0.71
1:A:300:THR:HG23	1:A:302:GLN:H	1.56	0.71
1:A:91:GLU:OE2	1:A:95:ARG:HG3	1.89	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
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
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:HE	1:A:310:ASP:CG	2.00	0.64
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: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
1:A:325:LEU:C	1:A:326:THR:O	2.35	0.61
2:B:86:ARG:CG	2:B:86:ARG:HH11	2.04	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: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:169:ASP:C	1:A:169:ASP:OD2	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:253:VAL:HG12	1:A:258:ARG:HH21	1.69	0.57
1:A:27:ILE:HG12	1:A:283:LEU:HD22	1.86	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
2:B:137:ASN:C	2:B:137:ASN:ND2	2.59	0.56
1:A:302:GLN:HE21	1:A:347:ASN:H	1.53	0.56
1:A:137:VAL:HG23	1:A:144:TYR:HE2	1.69	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:285:ARG:C	1:A:285:ARG:HD2	2.26	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:297:GLN:CA	1:A:297:GLN:HE21	2.21	0.52
1:A:50:LEU:HD21	1:A:119:ARG:HD2	1.91	0.52
2:B:116:LEU:CD1	2:B:124:MET:HE2	2.40	0.52
1:A:361:GLY:O	1:A:362:LEU:HB2	2.11	0.51
1:A:269:ILE:CG2	1:A:298:HIS:HA	2.41	0.51
1:A:304:ASN:HD22	1:A:305:PRO:HD2	1.75	0.51
1:A:197:HIS:HD2	1:A:199:SER:OG	1.94	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:CG	2:B:86:ARG:NH1	2.66	0.50
1:A:301:GLU:HG3	1:A:301:GLU:O	2.12	0.50
1:A:210:THR:HG21	1:A:213:ASP:H	1.76	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:223:ARG:C	1:A:223:ARG:HE	2.16	0.48
1:A:75:TYR:HB3	1:A:176:VAL:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:269:ILE:HG22	1:A:269:ILE:O	2.15	0.45
1:A:253:VAL:HG11	2:B:87:GLU:HB3	1.98	0.45
1:A:253:VAL:CG1	1:A:253:VAL:O	2.64	0.45
1:A:135:GLY:C	1:A:135:GLY:N	2.62	0.44
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:319:THR:OG1	1:A:321:GLU:HG3	2.19	0.43
1:A:35:ASN:HD22	1:A:318:ALA:HB1	1.83	0.43
1:A:258:ARG:HH22	2:B:87:GLU:CD	2.22	0.43
1:A:139:SER:HB3	1:A:144:TYR:CG	2.54	0.43
1:A:106:HIS:N	1:A:106:HIS:ND1	2.67	0.42
2:B:146:THR:O	2:B:146:THR:HG22	2.20	0.42
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:133:ALA:HB3	1:A:136:VAL:CA	2.46	0.41
1:A:132:MET:CE	1:A:135:GLY:N	2.83	0.41
1:A:41:ARG:HH21	1:A:310:ASP:CB	2.32	0.41
2:B:93:ASP:CG	2:B:97:ASN:O	2.59	0.41
1:A:75:TYR:HB3	1:A:176:VAL:CG2	2.50	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:302:GLN:HE21	1:A:347:ASN:N	2.17	0.41
1:A:196:PRO:HB3	1:A:275:PHE:CE2	2.56	0.41
1:A:331:LYS:HE2	1:A:331:LYS:HB3	1.79	0.41
1:A:133:ALA:CB	1:A:136:VAL:HG23	2.50	0.40
2:B:124:MET:HE3	2:B:124:MET:HB2	1.94	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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 (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	347/364 (95%)	327 (94%)	17 (5%)	3 (1%)	21 15
2	B	67/74 (90%)	64 (96%)	2 (3%)	1 (2%)	13 7
All	All	414/438 (94%)	391 (94%)	19 (5%)	4 (1%)	19 13

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 (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	271/280 (97%)	242 (89%)	29 (11%)	8 4
2	B	59/64 (92%)	52 (88%)	7 (12%)	6 3
All	All	330/344 (96%)	294 (89%)	36 (11%)	8 4

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	24	LEU
1	A	35	ASN
1	A	57	THR
1	A	84	LYS
1	A	95	ARG
1	A	117	LYS
1	A	119	ARG
1	A	125	GLN
1	A	140	ASN
1	A	151	VAL
1	A	163	TYR
1	A	166	LYS
1	A	169	ASP
1	A	170	ASP
1	A	176	VAL
1	A	204	SER
1	A	207	SER
1	A	221	ARG
1	A	223	ARG
1	A	244	ILE
1	A	285	ARG
1	A	290	VAL
1	A	297	GLN
1	A	300	THR
1	A	304	ASN
1	A	312	LYS
1	A	359	ASP
1	A	362	LEU
2	B	95	ASP
2	B	105	LEU
2	B	110	THR
2	B	122	ASP

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Mol	Chain	Res	Type
2	B	125	ILE
2	B	137	ASN
2	B	139	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) 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	297	GLN
1	A	298	HIS
1	A	302	GLN
1	A	304	ASN
1	A	336	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 carbohydrates in this entry.

5.6 Ligand geometry

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

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, 95^{th} 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%)	0.51	33 (9%) 11 14	13, 25, 46, 56	0
2	B	69/74 (93%)	0.16	2 (2%) 55 63	14, 23, 38, 49	0
All	All	420/438 (95%)	0.45	35 (8%) 14 19	13, 25, 46, 56	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	364	ALA	8.8
1	A	233	LEU	6.7
1	A	224	ARG	6.0
1	A	8	GLY	6.0
1	A	225	ALA	5.9
1	A	134	ASP	5.8
1	A	133	ALA	5.0
1	A	155	SER	5.0
1	A	363	GLY	4.9
1	A	167	GLY	3.9
1	A	168	GLY	3.6
1	A	156	ASP	3.6
1	A	63	HIS	3.5
1	A	310	ASP	3.2
1	A	309	ALA	3.2
1	A	95	ARG	3.2
1	A	89	ALA	3.1
1	A	135	GLY	3.0
1	A	194	ILE	2.9
2	B	79	THR	2.9
1	A	223	ARG	2.8
1	A	169	ASP	2.6
1	A	91	GLU	2.6
1	A	98	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	166	LYS	2.4
1	A	361	GLY	2.3
1	A	7	ALA	2.3
1	A	343	VAL	2.3
1	A	170	ASP	2.3
1	A	132	MET	2.2
1	A	96	ALA	2.1
1	A	92	VAL	2.1
1	A	221	ARG	2.1
1	A	359	ASP	2.0
2	B	139	GLU	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 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, 95^{th} 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
3	CA	B	800	1/1	1.00	0.07	-1.71	21,21,21,21	0
3	CA	B	801	1/1	0.99	0.04	-2.50	26,26,26,26	0

6.5 Other polymers (i)

There are no such residues in this entry.