



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:19 AM GMT

PDB ID : 2COL
Title : Crystal structure analysis of CyaA/C-Cam with Pyrophosphate
Authors : Guo, Q.; Tang, W.J.; Shen, Y.
Deposited on : 2005-05-18
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

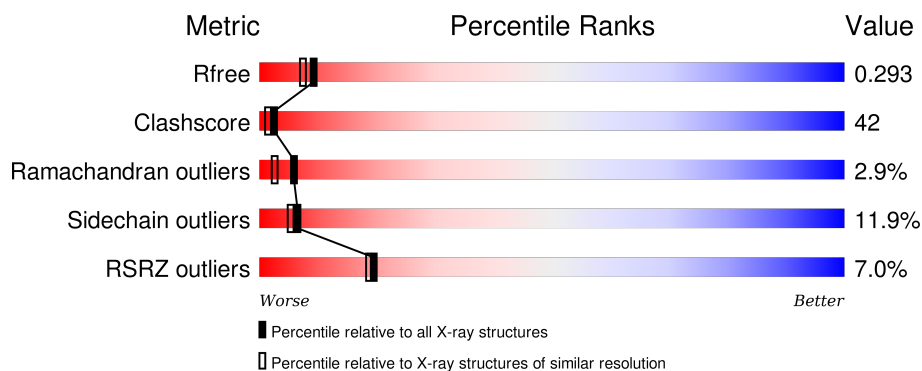
1 Overall quality at a glance

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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (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 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 $\leq 5\%$. 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	356	
2	B	67	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3345 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	349	Total	C	N	O	S	0	0	0
			2675	1659	493	517	6			

- Molecule 2 is a protein called Calmodulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	67	Total	C	N	O	S	0	0	0
			539	328	88	119	4			

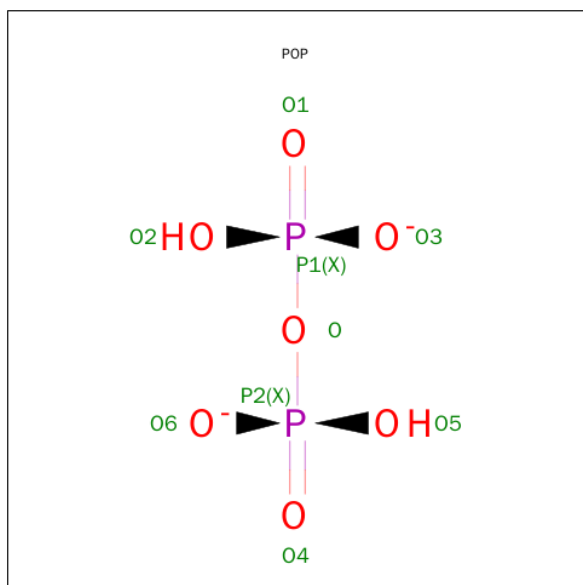
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: H₂O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			9	7	2		

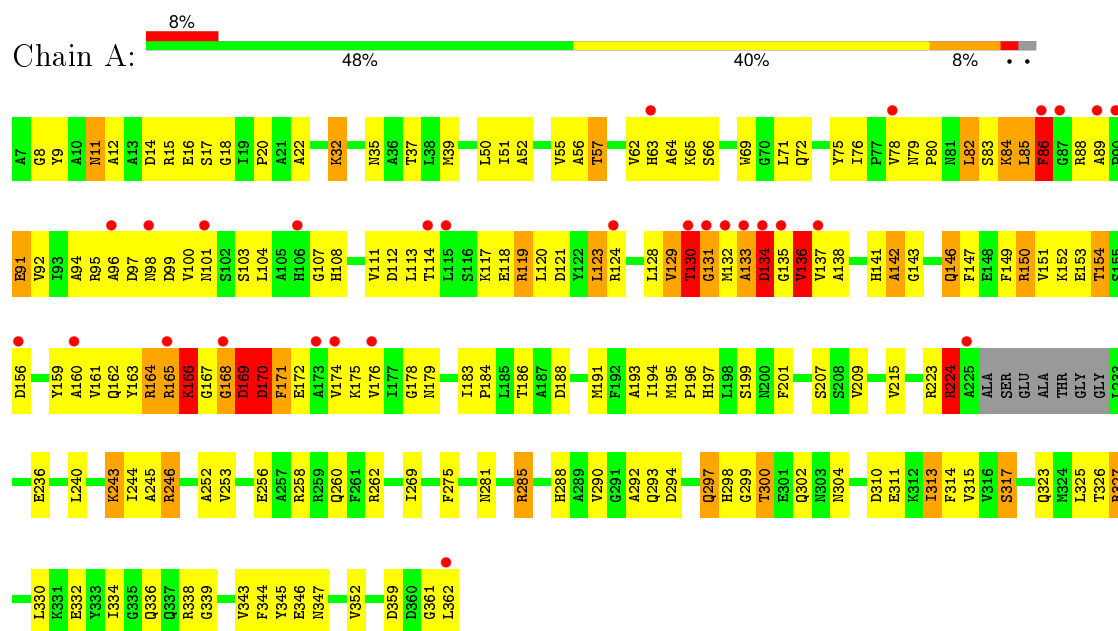
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	98	Total	O	0	0
			98	98		
6	B	20	Total	O	0	0
			20	20		

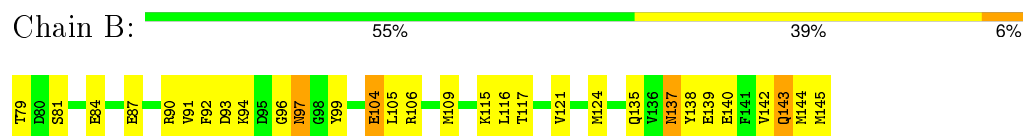
3 Residue-property plots [i](#)

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



- 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	79.61Å 79.61Å 139.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 43.80 – 2.17	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.20) 99.9 (43.80-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.22 (at 2.18Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.293 0.296 , 0.293	Depositor DCC
R_{free} test set	1168 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 24369 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3345	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 6.51% 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, 0.375 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, MG, POP

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	1.24	8/2719 (0.3%)	1.06	17/3670 (0.5%)
2	B	0.55	0/544	0.68	0/729
All	All	1.15	8/3263 (0.2%)	1.01	17/4399 (0.4%)

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	4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	130	THR	C-N	-34.15	0.71	1.33
1	A	86	PHE	C-N	-30.23	0.78	1.33
1	A	133	ALA	C-N	-18.30	0.92	1.34
1	A	131	GLY	N-CA	17.08	1.71	1.46
1	A	129	VAL	C-N	16.47	1.72	1.34
1	A	85	LEU	C-N	12.63	1.63	1.34
1	A	171	PHE	N-CA	9.39	1.65	1.46
1	A	166	LYS	C-N	5.66	1.43	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	PHE	O-C-N	-24.15	82.14	123.20
1	A	86	PHE	CA-C-N	17.59	151.39	116.20
1	A	86	PHE	C-N-CA	11.06	145.54	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	133	ALA	N-CA-C	-9.96	84.09	111.00
1	A	168	GLY	N-CA-C	-9.68	88.91	113.10
1	A	134	ASP	CA-C-N	-8.86	98.49	116.20
1	A	171	PHE	N-CA-C	8.78	134.71	111.00
1	A	131	GLY	N-CA-C	-8.44	92.00	113.10
1	A	130	THR	CA-C-N	-7.31	101.58	116.20
1	A	165	ARG	CA-C-N	-7.22	101.32	117.20
1	A	169	ASP	CA-C-N	-7.13	101.50	117.20
1	A	165	ARG	O-C-N	6.82	133.60	122.70
1	A	85	LEU	O-C-N	-6.63	112.09	122.70
1	A	171	PHE	N-CA-CB	-6.31	99.24	110.60
1	A	85	LEU	C-N-CA	6.06	136.84	121.70
1	A	166	LYS	C-N-CA	5.93	134.76	122.30
1	A	129	VAL	O-C-N	5.56	131.59	122.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	130	THR	Mainchain,Peptide
1	A	170	ASP	Peptide
1	A	86	PHE	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	2675	0	2612	244	0
2	B	539	0	490	28	0
3	B	2	0	0	0	0
4	A	2	0	0	0	0
5	A	9	0	0	2	0
6	A	98	0	0	5	0
6	B	20	0	0	2	0
All	All	3345	0	3102	265	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 42.

All (265) 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:131:GLY:N	1:A:131:GLY:CA	1.71	1.51
1:A:130:THR:CA	1:A:131:GLY:N	1.82	1.43
1:A:129:VAL:C	1:A:130:THR:N	1.71	1.41
1:A:124:ARG:NE	1:A:132:MET:CE	1.82	1.41
1:A:124:ARG:CZ	1:A:132:MET:HE1	1.59	1.30
1:A:124:ARG:NE	1:A:132:MET:HE1	1.41	1.23
1:A:86:PHE:CE1	1:A:143:GLY:HA2	1.72	1.23
1:A:130:THR:O	1:A:131:GLY:N	1.79	1.16
1:A:124:ARG:CZ	1:A:132:MET:CE	2.22	1.15
1:A:134:ASP:OD2	1:A:134:ASP:O	1.68	1.12
1:A:130:THR:C	1:A:131:GLY:CA	2.20	1.09
1:A:164:ARG:HH21	1:A:167:GLY:HA2	0.92	1.05
1:A:164:ARG:NH2	1:A:167:GLY:HA2	1.73	1.04
1:A:215:VAL:HG13	1:A:240:LEU:HD23	1.41	1.02
1:A:86:PHE:CZ	1:A:143:GLY:HA2	1.97	1.00
1:A:151:VAL:HG21	1:A:159:TYR:HB3	1.46	0.97
1:A:95:ARG:HH11	1:A:95:ARG:HB3	1.27	0.97
1:A:103:SER:O	1:A:108:HIS:HB2	1.67	0.94
1:A:119:ARG:HH11	1:A:119:ARG:HG2	1.34	0.92
1:A:297:GLN:HA	1:A:297:GLN:HE21	1.33	0.90
1:A:95:ARG:NH1	1:A:95:ARG:HB3	1.88	0.89
1:A:327:ARG:CB	1:A:327:ARG:HH11	1.85	0.88
1:A:75:TYR:HB3	1:A:176:VAL:HG21	1.55	0.88
1:A:153:GLU:HG2	1:A:154:THR:H	1.38	0.88
1:A:57:THR:HG23	1:A:188:ASP:HB3	1.56	0.87
1:A:96:ALA:O	1:A:100:VAL:HG23	1.74	0.86
1:A:50:LEU:HB3	1:A:55:VAL:HG21	1.58	0.86
1:A:124:ARG:CD	1:A:132:MET:HE1	2.05	0.85
1:A:79:ASN:HB3	1:A:82:LEU:HG	1.56	0.85
1:A:130:THR:CB	1:A:131:GLY:N	2.40	0.84
1:A:124:ARG:HE	1:A:132:MET:CE	1.85	0.84
1:A:302:GLN:HE21	1:A:346:GLU:HA	1.42	0.84
1:A:98:ASN:HA	1:A:101:ASN:ND2	1.91	0.83
1:A:300:THR:HG23	1:A:302:GLN:H	1.42	0.82
1:A:164:ARG:HD3	1:A:171:PHE:CZ	2.15	0.82
1:A:315:VAL:HG21	1:A:325:LEU:HD12	1.63	0.81
1:A:124:ARG:CZ	1:A:132:MET:HE3	2.11	0.81
1:A:131:GLY:C	1:A:131:GLY:N	2.34	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ARG:NE	1:A:132:MET:HE2	1.92	0.80
1:A:124:ARG:CD	1:A:132:MET:CE	2.59	0.80
1:A:327:ARG:CG	1:A:327:ARG:HH11	1.94	0.80
1:A:120:LEU:HG	1:A:124:ARG:HH11	1.47	0.78
1:A:151:VAL:CG2	1:A:159:TYR:HB3	2.13	0.78
1:A:133:ALA:O	1:A:135:GLY:N	2.17	0.76
1:A:130:THR:C	1:A:131:GLY:N	0.71	0.76
1:A:327:ARG:HB2	1:A:327:ARG:HH11	1.51	0.76
1:A:86:PHE:CD1	1:A:143:GLY:HA2	2.20	0.75
1:A:183:ILE:HG21	1:A:293:GLN:NE2	2.01	0.75
1:A:119:ARG:HH11	1:A:119:ARG:CG	2.00	0.74
2:B:116:LEU:HD13	2:B:124:MET:HE1	1.70	0.74
1:A:124:ARG:NE	1:A:132:MET:HE3	1.95	0.73
1:A:98:ASN:HA	1:A:101:ASN:HD22	1.50	0.73
2:B:137:ASN:HD22	2:B:137:ASN:C	1.91	0.73
1:A:362:LEU:HD23	2:B:90:ARG:HH22	1.54	0.72
1:A:151:VAL:HG23	1:A:160:ALA:O	1.90	0.70
1:A:153:GLU:HG2	1:A:154:THR:N	2.07	0.69
1:A:113:LEU:HG	1:A:161:VAL:HG21	1.73	0.68
1:A:12:ALA:O	1:A:16:GLU:HG3	1.93	0.68
2:B:137:ASN:ND2	2:B:140:GLU:H	1.92	0.68
1:A:302:GLN:NE2	1:A:346:GLU:HA	2.09	0.67
1:A:302:GLN:O	1:A:347:ASN:HA	1.95	0.67
2:B:138:TYR:O	2:B:142:VAL:HG13	1.95	0.66
1:A:288:HIS:HB2	6:A:981:HOH:O	1.94	0.66
1:A:269:ILE:HG22	1:A:269:ILE:O	1.96	0.66
1:A:124:ARG:NH2	1:A:132:MET:HE3	2.12	0.65
1:A:39:MET:HG2	1:A:315:VAL:HG22	1.77	0.65
1:A:82:LEU:N	1:A:82:LEU:HD23	2.10	0.65
1:A:179:ASN:HB2	1:A:183:ILE:O	1.96	0.65
1:A:11:ASN:ND2	1:A:14:ASP:H	1.95	0.65
1:A:95:ARG:HH11	1:A:95:ARG:CB	2.07	0.64
1:A:183:ILE:HG21	1:A:293:GLN:HE21	1.62	0.64
1:A:15:ARG:HH11	1:A:15:ARG:HG2	1.63	0.64
1:A:311:GLU:HG2	6:A:977:HOH:O	1.98	0.63
1:A:164:ARG:HD2	1:A:169:ASP:OD1	1.98	0.63
1:A:75:TYR:HB3	1:A:176:VAL:CG2	2.27	0.63
1:A:154:THR:C	1:A:156:ASP:H	2.02	0.63
2:B:92:PHE:O	2:B:104:GLU:HG2	1.99	0.63
1:A:57:THR:HG23	1:A:188:ASP:CB	2.28	0.63
1:A:111:VAL:HG12	1:A:112:ASP:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:THR:N	1:A:138:ALA:O	2.33	0.62
1:A:153:GLU:CG	1:A:154:THR:H	2.10	0.62
2:B:116:LEU:CD1	2:B:124:MET:HE1	2.30	0.61
1:A:223:ARG:HD3	1:A:236:GLU:OE2	2.00	0.61
1:A:269:ILE:HG23	1:A:298:HIS:HA	1.82	0.61
1:A:114:THR:HG22	1:A:160:ALA:HA	1.83	0.61
1:A:86:PHE:CZ	1:A:143:GLY:CA	2.79	0.61
1:A:129:VAL:CA	1:A:130:THR:N	2.64	0.60
1:A:64:ALA:O	1:A:76:ILE:HG23	2.01	0.60
1:A:300:THR:CG2	1:A:302:GLN:H	2.12	0.60
1:A:183:ILE:HD12	1:A:293:GLN:HE22	1.66	0.60
1:A:161:VAL:HG12	1:A:174:VAL:HG21	1.84	0.60
1:A:86:PHE:CE1	1:A:143:GLY:CA	2.66	0.59
1:A:338:ARG:O	2:B:94:LYS:HG2	2.02	0.59
1:A:201:PHE:CZ	1:A:252:ALA:HA	2.37	0.59
2:B:109:MET:HG3	2:B:124:MET:HE1	1.83	0.59
1:A:197:HIS:CD2	1:A:199:SER:H	2.19	0.59
1:A:362:LEU:CD2	2:B:90:ARG:HH22	2.16	0.58
1:A:32:LYS:NZ	1:A:32:LYS:HB2	2.18	0.58
1:A:50:LEU:HB3	1:A:55:VAL:CG2	2.30	0.58
1:A:91:GLU:O	1:A:95:ARG:HG3	2.04	0.58
1:A:95:ARG:HA	1:A:98:ASN:ND2	2.18	0.58
1:A:118:GLU:O	1:A:121:ASP:HB2	2.03	0.58
1:A:79:ASN:HB3	1:A:82:LEU:CG	2.33	0.57
1:A:311:GLU:O	1:A:326:THR:HA	2.04	0.57
2:B:99:TYR:CD1	2:B:135:GLN:NE2	2.72	0.57
1:A:332:GLU:HG2	1:A:336:GLN:HE21	1.70	0.57
1:A:338:ARG:HD2	6:B:814:HOH:O	2.03	0.57
1:A:258:ARG:HH22	2:B:87:GLU:CD	2.08	0.57
1:A:113:LEU:HG	1:A:161:VAL:CG2	2.33	0.57
1:A:132:MET:HA	1:A:136:VAL:O	2.05	0.57
1:A:327:ARG:HH11	1:A:327:ARG:HG3	1.69	0.56
1:A:195:MET:HG2	1:A:344:PHE:HB2	1.87	0.56
1:A:124:ARG:HD2	1:A:132:MET:HE1	1.87	0.56
1:A:327:ARG:CG	1:A:327:ARG:NH1	2.61	0.56
2:B:116:LEU:HD13	2:B:124:MET:CE	2.34	0.55
2:B:81:SER:N	2:B:84:GLU:OE1	2.35	0.55
1:A:169:ASP:CG	1:A:170:ASP:H	2.05	0.55
1:A:244:ILE:HD12	1:A:244:ILE:C	2.27	0.55
1:A:133:ALA:O	1:A:134:ASP:C	2.36	0.54
1:A:285:ARG:HD2	1:A:285:ARG:C	2.26	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:117:THR:O	2:B:121:VAL:HG23	2.07	0.54
1:A:95:ARG:HA	1:A:98:ASN:HD22	1.72	0.54
1:A:327:ARG:HB2	1:A:327:ARG:NH1	2.18	0.54
1:A:281:ASN:O	1:A:285:ARG:HB2	2.08	0.54
1:A:130:THR:OG1	1:A:138:ALA:O	2.25	0.54
1:A:183:ILE:CD1	1:A:293:GLN:HE22	2.20	0.54
1:A:141:HIS:O	1:A:143:GLY:N	2.41	0.53
1:A:327:ARG:NH1	1:A:327:ARG:HG3	2.23	0.53
1:A:339:GLY:HA2	2:B:94:LYS:HE3	1.90	0.53
1:A:71:LEU:HD13	1:A:161:VAL:HG13	1.90	0.53
1:A:123:LEU:HD13	1:A:128:LEU:HB2	1.91	0.53
1:A:94:ALA:HA	1:A:97:ASP:OD2	2.09	0.53
1:A:197:HIS:HD2	1:A:199:SER:H	1.56	0.53
1:A:330:LEU:O	1:A:334:ILE:HG13	2.09	0.52
1:A:209:VAL:HB	1:A:244:ILE:HG22	1.92	0.52
1:A:246:ARG:HH11	1:A:246:ARG:HB3	1.73	0.52
2:B:137:ASN:HD21	2:B:140:GLU:H	1.55	0.52
1:A:124:ARG:HG3	1:A:129:VAL:HG23	1.90	0.52
1:A:253:VAL:HG13	1:A:345:TYR:CD2	2.44	0.52
1:A:224:ARG:CD	1:A:224:ARG:H	2.23	0.52
1:A:298:HIS:HD2	1:A:299:GLY:O	1.93	0.51
1:A:119:ARG:NH1	1:A:119:ARG:CG	2.64	0.51
1:A:65:LYS:HZ3	5:A:893:POP:P2	2.33	0.51
1:A:50:LEU:O	1:A:55:VAL:HG23	2.11	0.51
1:A:171:PHE:CD2	1:A:171:PHE:N	2.77	0.51
1:A:15:ARG:NH1	1:A:15:ARG:HG2	2.24	0.51
1:A:57:THR:HB	1:A:294:ASP:O	2.11	0.51
1:A:130:THR:O	1:A:131:GLY:CA	2.59	0.51
1:A:269:ILE:HD12	1:A:297:GLN:O	2.10	0.51
1:A:315:VAL:CG2	1:A:325:LEU:HD12	2.38	0.50
1:A:260:GLN:HE21	1:A:262:ARG:HD2	1.76	0.50
1:A:124:ARG:CD	1:A:132:MET:HE2	2.34	0.50
1:A:146:GLN:O	1:A:166:LYS:HB2	2.11	0.50
1:A:65:LYS:NZ	5:A:893:POP:O4	2.43	0.50
1:A:17:SER:HB2	6:A:972:HOH:O	2.12	0.50
1:A:78:VAL:HG23	1:A:175:LYS:C	2.32	0.50
1:A:132:MET:HB3	1:A:137:VAL:HA	1.94	0.50
1:A:297:GLN:HA	1:A:297:GLN:NE2	2.14	0.50
1:A:253:VAL:HG12	1:A:253:VAL:O	2.11	0.50
1:A:85:LEU:HD22	1:A:88:ARG:HG3	1.93	0.50
2:B:137:ASN:ND2	2:B:137:ASN:C	2.63	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ASN:C	1:A:11:ASN:HD22	2.15	0.49
1:A:69:TRP:HB3	1:A:141:HIS:CD2	2.47	0.49
1:A:103:SER:HB3	1:A:108:HIS:CD2	2.47	0.49
1:A:132:MET:SD	1:A:137:VAL:HG12	2.51	0.49
2:B:139:GLU:O	2:B:142:VAL:HG22	2.11	0.49
1:A:20:PRO:HG2	1:A:51:ILE:HG22	1.95	0.49
1:A:78:VAL:O	1:A:80:PRO:HD3	2.12	0.49
1:A:114:THR:HA	1:A:159:TYR:O	2.13	0.49
1:A:57:THR:CG2	1:A:297:GLN:HG2	2.43	0.49
1:A:151:VAL:HG22	1:A:152:LYS:N	2.28	0.49
1:A:62:VAL:HG22	1:A:108:HIS:CD2	2.47	0.49
1:A:196:PRO:HB3	1:A:275:PHE:CE2	2.48	0.49
1:A:194:ILE:O	1:A:196:PRO:HD3	2.13	0.48
1:A:201:PHE:CE2	1:A:252:ALA:HA	2.48	0.48
1:A:78:VAL:HG13	1:A:100:VAL:CG1	2.43	0.48
1:A:83:SER:O	1:A:85:LEU:N	2.43	0.48
1:A:183:ILE:CD1	1:A:293:GLN:NE2	2.76	0.48
2:B:94:LYS:HD3	6:B:819:HOH:O	2.14	0.48
1:A:72:GLN:NE2	1:A:82:LEU:HB3	2.28	0.48
1:A:50:LEU:CB	1:A:55:VAL:HG21	2.39	0.47
1:A:163:TYR:CA	1:A:171:PHE:HD1	2.27	0.47
1:A:111:VAL:CG1	1:A:112:ASP:N	2.77	0.47
1:A:22:ALA:HB1	1:A:290:VAL:HB	1.97	0.47
1:A:120:LEU:HD22	1:A:159:TYR:CD1	2.51	0.46
2:B:144:MET:HG2	2:B:144:MET:O	2.14	0.46
1:A:163:TYR:C	1:A:171:PHE:HD1	2.18	0.46
1:A:117:LYS:O	1:A:120:LEU:HB3	2.15	0.46
1:A:124:ARG:NH1	1:A:132:MET:HE1	2.18	0.46
1:A:240:LEU:O	1:A:243:LYS:HB2	2.14	0.46
1:A:8:GLY:O	1:A:9:TYR:C	2.52	0.46
1:A:298:HIS:CD2	1:A:299:GLY:O	2.68	0.46
1:A:179:ASN:HB3	1:A:183:ILE:H	1.80	0.46
1:A:285:ARG:HD3	6:A:969:HOH:O	2.14	0.46
1:A:132:MET:HB3	1:A:137:VAL:HG12	1.98	0.46
1:A:197:HIS:HD2	1:A:199:SER:OG	1.98	0.46
1:A:20:PRO:HB3	1:A:52:ALA:HB2	1.98	0.46
1:A:343:VAL:HG22	2:B:91:VAL:HG13	1.98	0.46
1:A:164:ARG:HD2	1:A:169:ASP:HA	1.98	0.46
1:A:193:ALA:HA	1:A:300:THR:HG21	1.99	0.45
2:B:93:ASP:OD1	2:B:97:ASN:O	2.35	0.45
1:A:130:THR:C	1:A:131:GLY:C	2.76	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:HIS:O	1:A:142:ALA:C	2.55	0.45
1:A:178:GLY:HA2	1:A:184:PRO:HA	1.98	0.45
1:A:315:VAL:HG21	1:A:325:LEU:CD1	2.40	0.45
1:A:89:ALA:O	1:A:92:VAL:N	2.49	0.45
1:A:150:ARG:HG3	1:A:162:GLN:HB2	1.99	0.44
1:A:313:ILE:HD11	1:A:315:VAL:HG22	1.99	0.44
1:A:253:VAL:HG11	2:B:87:GLU:HB3	1.99	0.44
1:A:97:ASP:O	1:A:101:ASN:CG	2.56	0.44
1:A:179:ASN:HB3	1:A:183:ILE:N	2.32	0.44
2:B:109:MET:HG3	2:B:124:MET:CE	2.45	0.44
1:A:88:ARG:HH21	1:A:92:VAL:HG21	1.83	0.44
1:A:66:SER:O	1:A:84:LYS:HG2	2.17	0.44
1:A:104:LEU:HD23	1:A:108:HIS:O	2.18	0.44
1:A:72:GLN:HA	1:A:113:LEU:HD21	1.99	0.44
1:A:346:GLU:HG3	1:A:347:ASN:O	2.18	0.44
1:A:83:SER:C	1:A:85:LEU:H	2.19	0.44
1:A:150:ARG:NH1	1:A:150:ARG:HG2	2.33	0.44
1:A:76:ILE:C	1:A:176:VAL:HG23	2.38	0.44
2:B:106:ARG:HG2	2:B:106:ARG:HH11	1.83	0.44
1:A:94:ALA:O	1:A:97:ASP:HB2	2.17	0.43
1:A:98:ASN:CA	1:A:101:ASN:ND2	2.73	0.43
1:A:64:ALA:O	1:A:65:LYS:C	2.57	0.43
1:A:85:LEU:CD2	1:A:88:ARG:HG3	2.48	0.43
1:A:56:ALA:O	1:A:186:THR:HG22	2.18	0.43
1:A:297:GLN:CA	1:A:297:GLN:HE21	2.09	0.43
1:A:98:ASN:O	1:A:99:ASP:C	2.57	0.43
1:A:269:ILE:CG2	1:A:269:ILE:O	2.66	0.43
1:A:71:LEU:CD1	1:A:161:VAL:HG13	2.48	0.43
1:A:334:ILE:O	1:A:338:ARG:HG3	2.18	0.43
1:A:154:THR:C	1:A:156:ASP:N	2.70	0.43
1:A:244:ILE:HD12	1:A:245:ALA:N	2.34	0.43
1:A:313:ILE:HD12	1:A:314:PHE:N	2.33	0.42
1:A:361:GLY:O	1:A:362:LEU:HB2	2.19	0.42
1:A:104:LEU:HD23	1:A:104:LEU:HA	1.87	0.42
1:A:124:ARG:HE	1:A:132:MET:HE3	1.72	0.42
1:A:164:ARG:HG2	1:A:171:PHE:CE1	2.55	0.42
1:A:85:LEU:HD22	1:A:92:VAL:HG11	2.01	0.42
1:A:62:VAL:O	1:A:63:HIS:ND1	2.53	0.42
1:A:130:THR:HG1	1:A:138:ALA:C	2.22	0.42
1:A:79:ASN:O	1:A:82:LEU:HG	2.20	0.42
1:A:285:ARG:NH1	1:A:285:ARG:HG3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:GLY:O	1:A:137:VAL:HG13	2.19	0.41
1:A:18:GLY:N	6:A:925:HOH:O	2.52	0.41
1:A:162:GLN:HA	1:A:172:GLU:O	2.21	0.41
1:A:313:ILE:HD11	1:A:315:VAL:CG2	2.50	0.41
1:A:123:LEU:HD12	1:A:129:VAL:HG13	2.03	0.41
1:A:22:ALA:HB1	1:A:290:VAL:CB	2.51	0.41
1:A:128:LEU:N	1:A:128:LEU:CD1	2.84	0.41
1:A:201:PHE:CE1	1:A:252:ALA:HA	2.56	0.41
1:A:39:MET:HB3	1:A:313:ILE:CD1	2.51	0.41
1:A:11:ASN:O	1:A:15:ARG:HD2	2.20	0.41
2:B:143:GLN:C	2:B:145:MET:N	2.74	0.41
1:A:288:HIS:HA	1:A:292:ALA:O	2.21	0.40
1:A:88:ARG:HE	1:A:88:ARG:HB3	1.66	0.40
2:B:106:ARG:HB3	2:B:106:ARG:CZ	2.50	0.40
1:A:193:ALA:HB2	1:A:302:GLN:HG2	2.03	0.40
1:A:300:THR:CG2	1:A:302:GLN:HB2	2.51	0.40
1:A:246:ARG:HA	1:A:246:ARG:HD2	1.74	0.40
1:A:191:MET:HB3	1:A:191:MET:HE2	1.99	0.40
1:A:130:THR:OG1	1:A:138:ALA:HB3	2.19	0.40
1:A:103:SER:O	1:A:108:HIS:CB	2.55	0.40
1:A:11:ASN:HD21	1:A:14:ASP:H	1.67	0.40
1:A:147:PHE:HB2	1:A:149:PHE:CE1	2.57	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	345/356 (97%)	301 (87%)	34 (10%)	10 (3%)	6	2
2	B	65/67 (97%)	60 (92%)	3 (5%)	2 (3%)	5	2
All	All	410/423 (97%)	361 (88%)	37 (9%)	12 (3%)	6	2

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	142	ALA
1	A	169	ASP
1	A	84	LYS
1	A	166	LYS
2	B	97	ASN
1	A	224	ARG
1	A	134	ASP
1	A	317	SER
2	B	96	GLY
1	A	136	VAL
1	A	107	GLY
1	A	168	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	271/274 (99%)	238 (88%)	33 (12%)	6	5
2	B	58/58 (100%)	52 (90%)	6 (10%)	9	8
All	All	329/332 (99%)	290 (88%)	39 (12%)	6	5

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	32	LYS
1	A	35	ASN
1	A	37	THR
1	A	57	THR
1	A	82	LEU
1	A	91	GLU
1	A	119	ARG
1	A	123	LEU
1	A	136	VAL
1	A	146	GLN

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Mol	Chain	Res	Type
1	A	150	ARG
1	A	154	THR
1	A	164	ARG
1	A	165	ARG
1	A	169	ASP
1	A	170	ASP
1	A	207	SER
1	A	224	ARG
1	A	243	LYS
1	A	246	ARG
1	A	256	GLU
1	A	285	ARG
1	A	297	GLN
1	A	300	THR
1	A	304	ASN
1	A	310	ASP
1	A	313	ILE
1	A	317	SER
1	A	323	GLN
1	A	327	ARG
1	A	352	VAL
1	A	359	ASP
2	B	79	THR
2	B	104	GLU
2	B	105	LEU
2	B	115	LYS
2	B	137	ASN
2	B	143	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	35	ASN
1	A	98	ASN
1	A	101	ASN
1	A	125	GLN
1	A	162	GLN
1	A	197	HIS
1	A	260	GLN
1	A	293	GLN
1	A	297	GLN

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Mol	Chain	Res	Type
1	A	298	HIS
1	A	302	GLN
1	A	304	ASN
1	A	336	GLN
2	B	111	ASN
2	B	135	GLN
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 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	POP	A	893	-	8,8,8	1.00	0	13,13,13	1.24	2 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	POP	A	893	-	-	0/6/6/6	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	893	POP	O5-P2-O4	-2.87	100.32	110.85
5	A	893	POP	P2-O-P1	2.05	138.49	132.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	893	POP	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

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	349/356 (98%)	0.57	29 (8%) 14 13	7, 26, 47, 56	0
2	B	67/67 (100%)	0.20	0 100 100	11, 22, 35, 47	0
All	All	416/423 (98%)	0.51	29 (6%) 19 19	7, 25, 47, 56	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	GLY	8.5
1	A	132	MET	5.4
1	A	133	ALA	4.9
1	A	173	ALA	4.6
1	A	168	GLY	4.3
1	A	130	THR	3.7
1	A	134	ASP	3.6
1	A	135	GLY	3.5
1	A	165	ARG	3.2
1	A	115	LEU	3.2
1	A	89	ALA	3.1
1	A	174	VAL	3.0
1	A	78	VAL	3.0
1	A	137	VAL	2.8
1	A	362	LEU	2.7
1	A	176	VAL	2.7
1	A	225	ALA	2.5
1	A	98	ASN	2.5
1	A	87	GLY	2.4
1	A	63	HIS	2.4
1	A	106	HIS	2.3
1	A	86	PHE	2.2
1	A	101	ASN	2.2
1	A	156	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	124	ARG	2.1
1	A	114	THR	2.1
1	A	96	ALA	2.1
1	A	160	ALA	2.0
1	A	90	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 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, 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	LLDF	B-factors(Å ²)	Q < 0.9
5	POP	A	893	9/9	0.71	0.25	1.74	71,72,72,73	0
4	MG	A	908	1/1	0.74	0.15	-0.27	23,23,23,23	0
3	CA	B	801	1/1	0.81	0.12	-1.46	33,33,33,33	0
3	CA	B	800	1/1	0.96	0.08	-2.54	24,24,24,24	0
4	MG	A	907	1/1	0.79	0.28	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.