



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:42 PM UTC

PDB ID : 2COL / pdb_00002col
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

<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

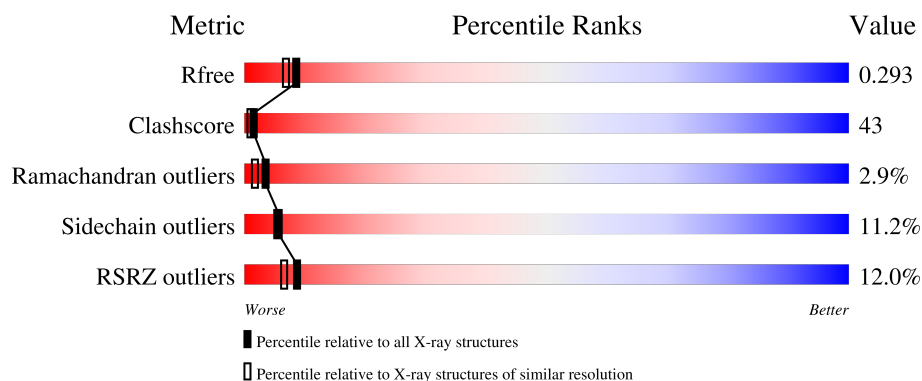
1 Overall quality at a glance

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	356	
2	B	67	

2 Entry composition [i](#)

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			

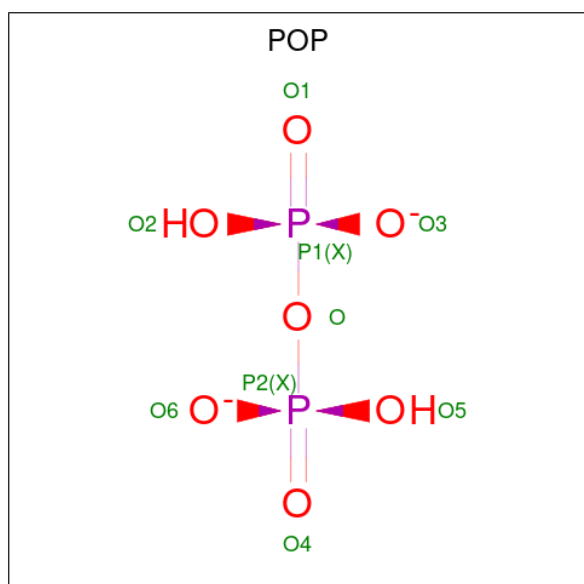
- Molecule 2 is a protein called Calmodulin.

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 MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).



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

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

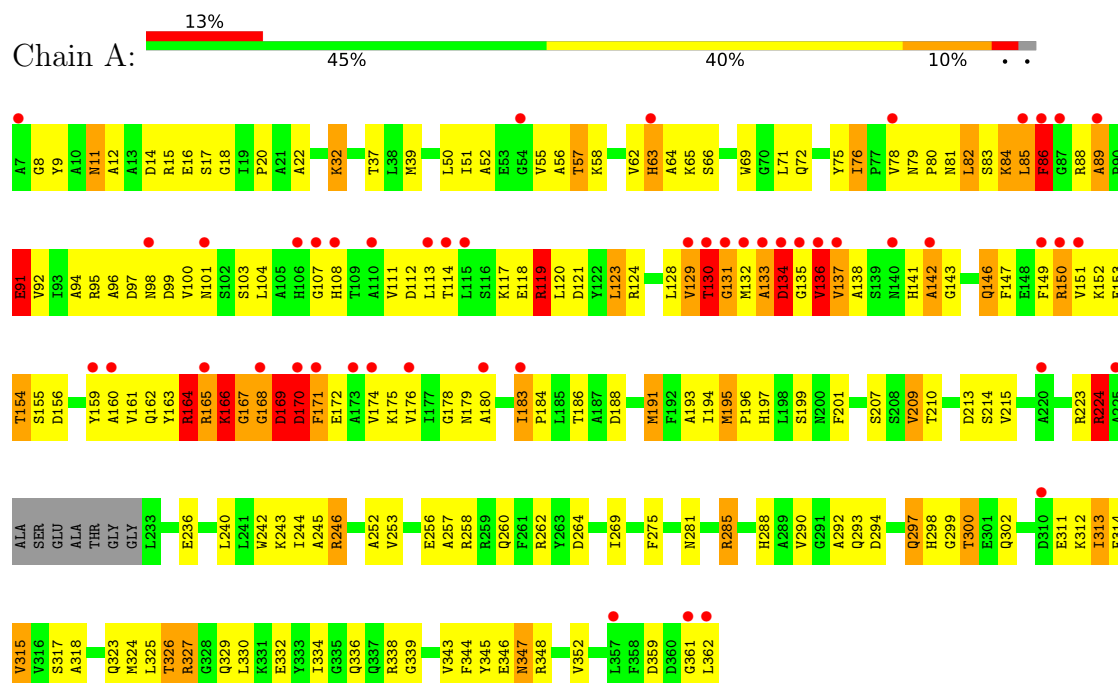
- Molecule 6 is water.

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

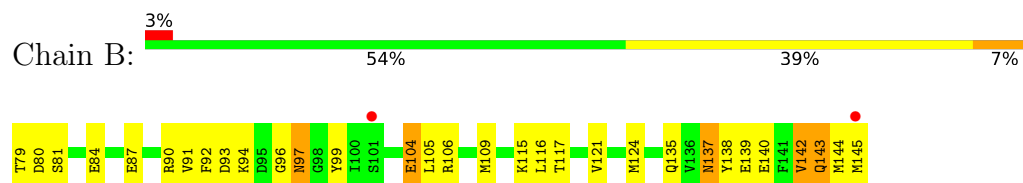
3 Residue-property plots

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	79.61Å 79.61Å 139.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.20) 99.9 (50.00-2.20)	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	1220 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.5	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	0.88	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, 0.333 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: MG, CA, 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	1.65	12/2719 (0.4%)	1.45	52/3670 (1.4%)
2	B	0.69	0/544	1.01	1/729 (0.1%)
All	All	1.53	12/3263 (0.4%)	1.38	53/4399 (1.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	130	THR	C-N	-42.41	0.71	1.33
1	A	86	PHE	C-N	-37.58	0.78	1.33
1	A	133	ALA	C-N	-29.80	0.92	1.33
1	A	129	VAL	C-N	27.31	1.72	1.33
1	A	85	LEU	C-N	21.00	1.63	1.33
1	A	131	GLY	N-CA	18.13	1.71	1.45
1	A	171	PHE	N-CA	14.69	1.65	1.46
1	A	170	ASP	N-CA	-7.25	1.36	1.46
1	A	166	LYS	C-N	6.66	1.43	1.33
1	A	167	GLY	C-N	-5.54	1.25	1.33
1	A	168	GLY	C-N	-5.49	1.26	1.33
1	A	191	MET	SD-CE	-5.18	1.66	1.79

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	PHE	O-C-N	-30.41	82.14	122.59
1	A	133	ALA	N-CA-C	-16.52	84.09	110.20
1	A	86	PHE	CA-C-N	12.31	145.54	121.41
1	A	86	PHE	C-N-CA	12.31	145.54	121.41
1	A	171	PHE	N-CA-C	11.23	134.71	110.80
1	A	180	ALA	N-CA-C	-10.79	98.63	112.23
1	A	168	GLY	N-CA-C	-10.24	88.91	113.18
1	A	165	ARG	O-C-N	9.11	133.60	123.22
1	A	131	GLY	N-CA-C	-8.94	92.00	113.18
1	A	91	GLU	N-CA-C	-8.26	102.22	111.14
1	A	85	LEU	CA-C-N	8.01	136.84	121.54
1	A	85	LEU	C-N-CA	8.01	136.84	121.54
1	A	85	LEU	O-C-N	-7.99	112.09	121.99
1	A	134	ASP	CA-CB-CG	-7.93	104.67	112.60
1	A	129	VAL	O-C-N	7.54	131.59	123.00
1	A	214	SER	N-CA-C	-7.46	99.98	110.50
1	A	315	VAL	N-CA-C	7.03	118.60	108.48
1	A	209	VAL	N-CA-C	6.91	117.50	111.56
1	A	166	LYS	CA-C-N	6.81	134.76	121.41
1	A	166	LYS	C-N-CA	6.81	134.76	121.41
1	A	155	SER	N-CA-C	-6.81	103.94	113.20
1	A	171	PHE	N-CA-CB	-6.66	99.24	110.49
1	A	195	MET	CA-C-N	6.46	126.22	119.76
1	A	195	MET	C-N-CA	6.46	126.22	119.76
1	A	164	ARG	O-C-N	-6.39	115.97	123.25
1	A	133	ALA	O-C-N	6.36	130.51	123.01
1	A	165	ARG	CA-C-N	6.30	133.58	121.54
1	A	165	ARG	C-N-CA	6.30	133.58	121.54
1	A	76	ILE	CA-C-N	6.17	126.12	119.76
1	A	76	ILE	C-N-CA	6.17	126.12	119.76
1	A	119	ARG	N-CA-C	-6.14	104.67	111.36
1	A	326	THR	N-CA-C	-6.12	102.28	110.55
1	A	347	ASN	N-CA-C	-6.12	101.61	110.24
1	A	170	ASP	CA-C-N	-6.11	109.87	121.54
1	A	170	ASP	C-N-CA	-6.11	109.87	121.54
1	A	134	ASP	O-C-N	6.11	130.72	122.59
1	A	348	ARG	N-CA-C	6.01	120.23	113.01
1	A	318	ALA	N-CA-C	-5.71	106.39	113.19
1	A	169	ASP	O-C-N	5.70	130.17	122.59
1	A	169	ASP	N-CA-C	5.61	122.74	110.80
1	A	89	ALA	N-CA-C	5.45	118.55	110.32
1	A	169	ASP	CA-C-O	5.43	128.28	120.51
1	A	133	ALA	CA-C-N	-5.36	111.31	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	133	ALA	C-N-CA	-5.36	111.31	121.54
1	A	170	ASP	CA-CB-CG	-5.36	107.24	112.60
1	A	264	ASP	N-CA-C	-5.35	99.06	108.20
1	A	58	LYS	N-CA-C	5.33	118.91	109.96
1	A	183	ILE	CA-C-N	5.33	125.32	119.89
1	A	183	ILE	C-N-CA	5.33	125.32	119.89
1	A	130	THR	O-C-N	5.32	129.66	122.59
1	A	257	ALA	N-CA-C	5.19	118.71	112.38
2	B	80	ASP	N-CA-C	-5.18	103.56	110.55
1	A	329	GLN	N-CA-C	-5.08	106.59	112.89

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	252	0
2	B	539	0	490	30	0
3	A	2	0	0	0	0
4	A	9	0	0	2	0
5	B	2	0	0	0	0
6	A	98	0	0	5	0
6	B	20	0	0	2	0
All	All	3345	0	3102	272	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 43.

All (272) 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:129:VAL:C	1:A:130:THR:N	1.71	1.46
1:A:130:THR:CA	1:A:131:GLY:N	1.82	1.43
1:A:124:ARG:NE	1:A:132:MET:CE	1.82	1.41
1:A:124:ARG:CZ	1:A:132:MET:HE3	1.62	1.29
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:124:ARG:CD	1:A:132:MET:HE1	1.77	1.13
1:A:130:THR:C	1:A:131:GLY:CA	2.20	1.13
1:A:134:ASP:O	1:A:134:ASP:OD2	1.68	1.12
1:A:124:ARG:NE	1:A:132:MET:HE1	1.48	1.12
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:95:ARG:HH11	1:A:95:ARG:HB3	1.27	0.97
1:A:151:VAL:HG21	1:A:159:TYR:HB3	1.46	0.97
1:A:124:ARG:NH2	1:A:132:MET:HE3	1.78	0.97
1:A:103:SER:O	1:A:108:HIS:HB2	1.67	0.94
1:A:119:ARG:HG2	1:A:119:ARG:HH11	1.34	0.92
2:B:109:MET:HG3	2:B:124:MET:HE1	1.50	0.91
1:A:297:GLN:HE21	1:A:297:GLN:HA	1.33	0.90
1:A:95:ARG:HB3	1:A:95:ARG:NH1	1.88	0.89
1:A:124:ARG:HD2	1:A:132:MET:HE1	1.55	0.88
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:131:GLY:N	1:A:131:GLY:C	2.34	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:130:THR:C	1:A:131:GLY:N	0.71	0.81
1:A:327:ARG:HH11	1:A:327:ARG:CG	1.94	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:B:137:ASN:C	2:B:137:ASN:HD22	1.91	0.78
1:A:133:ALA:O	1:A:135:GLY:N	2.17	0.76
1:A:327:ARG:HH11	1:A:327:ARG:HB2	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
1:A:98:ASN:HA	1:A:101:ASN:HD22	1.50	0.73
1:A:124:ARG:NE	1:A:132:MET:HE3	1.75	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
2:B:116:LEU:HD13	2:B:124:MET:HE2	1.74	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
1:A:154:THR:C	1:A:156:ASP:H	2.02	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:124:ARG:CZ	1:A:132:MET:HE1	2.08	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: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:242:TRP:CD2	2:B:144:MET:HE1	2.32	0.64
1:A:183:ILE:HG21	1:A:293:GLN:HE21	1.62	0.64
1:A:15:ARG:HG2	1:A:15:ARG:HH11	1.63	0.64
1:A:311:GLU:HG2	6:A:977:HOH:O	1.98	0.63
1:A:75:TYR:HB3	1:A:176:VAL:CG2	2.27	0.63
1:A:164:ARG:HD2	1:A:169:ASP:OD1	1.98	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
1:A:258:ARG:HH22	2:B:87:GLU:CD	2.08	0.62
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ARG:HD3	1:A:236:GLU:OE2	2.00	0.61
1:A:124:ARG:NE	1:A:132:MET:HE2	2.06	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:285:ARG:HD2	1:A:285:ARG:C	2.26	0.59
1:A:244:ILE:HD12	1:A:244:ILE:C	2.27	0.59
1:A:338:ARG:O	2:B:94:LYS:HG2	2.02	0.59
1:A:133:ALA:O	1:A:134:ASP:C	2.36	0.59
1:A:201:PHE:CZ	1:A:252:ALA:HA	2.37	0.59
1:A:312:LYS:HB3	1:A:324:MET:HE3	1.85	0.59
1:A:197:HIS:CD2	1:A:199:SER:H	2.19	0.59
1:A:169:ASP:CG	1:A:170:ASP:H	2.05	0.58
1:A:32:LYS:HB2	1:A:32:LYS:NZ	2.18	0.58
1:A:362:LEU:CD2	2:B:90:ARG:HH22	2.16	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: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:242:TRP:CE3	2:B:144:MET:HE1	2.41	0.56
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
2:B:116:LEU:HD13	2:B:124:MET:CE	2.34	0.56
1:A:327:ARG:CG	1:A:327:ARG:NH1	2.61	0.56
2:B:81:SER:N	2:B:84:GLU:OE1	2.35	0.55
1:A:78:VAL:HG23	1:A:175:LYS:C	2.32	0.54
1:A:130:THR:C	1:A:131:GLY:C	2.76	0.54
2:B:117:THR:O	2:B:121:VAL:HG23	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ASN:HD22	1:A:11:ASN:C	2.15	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
2:B:137:ASN:C	2:B:137:ASN:ND2	2.63	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:134:ASP:OD2	1:A:134:ASP:C	2.39	0.54
1:A:183:ILE:CD1	1:A:293:GLN:HE22	2.20	0.54
1:A:124:ARG:CD	1:A:132:MET:CE	2.59	0.53
1:A:141:HIS:O	1:A:143:GLY:N	2.41	0.53
1:A:327:ARG:HG3	1:A:327:ARG:NH1	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:94:ALA:HA	1:A:97:ASP:OD2	2.09	0.53
1:A:123:LEU:HD13	1:A:128:LEU:HB2	1.91	0.53
1:A:197:HIS:HD2	1:A:199:SER:H	1.56	0.53
1:A:209:VAL:HB	1:A:244:ILE:HG22	1.92	0.52
1:A:330:LEU:O	1:A:334:ILE:HG13	2.09	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:224:ARG:CD	1:A:224:ARG:H	2.23	0.52
1:A:253:VAL:HG13	1:A:345:TYR:CD2	2.44	0.52
1:A:298:HIS:HD2	1:A:299:GLY:O	1.93	0.51
1:A:119:ARG:CG	1:A:119:ARG:NH1	2.64	0.51
1:A:50:LEU:O	1:A:55:VAL:HG23	2.11	0.51
1:A:15:ARG:HG2	1:A:15:ARG:NH1	2.24	0.51
1:A:171:PHE:CD2	1:A:171:PHE:N	2.77	0.51
1:A:163:TYR:C	1:A:171:PHE:HD1	2.18	0.51
1:A:8:GLY:O	1:A:9:TYR:C	2.52	0.51
1:A:57:THR:HB	1:A:294:ASP:O	2.11	0.51
1:A:269:ILE:HD12	1:A:297:GLN:O	2.10	0.51
1:A:65:LYS:NZ	4:A:893:POP:O4	2.43	0.50
1:A:260:GLN:HE21	1:A:262:ARG:HD2	1.76	0.50
1:A:315:VAL:CG2	1:A:325:LEU:HD12	2.38	0.50
1:A:146:GLN:O	1:A:166:LYS:HB2	2.11	0.50
1:A:17:SER:HB2	6:A:972:HOH:O	2.12	0.50
1:A:124:ARG:HE	1:A:132:MET:HE2	1.68	0.50
1:A:132:MET:HB3	1:A:137:VAL:HA	1.94	0.50
1:A:83:SER:C	1:A:85:LEU:H	2.19	0.50
1:A:85:LEU:HD22	1:A:88:ARG:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:VAL:HG12	1:A:253:VAL:O	2.11	0.50
1:A:297:GLN:HA	1:A:297:GLN:NE2	2.14	0.50
1:A:141:HIS:O	1:A:142:ALA:C	2.55	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:244:ILE:HD11	2:B:109:MET:HE1	1.95	0.49
1:A:57:THR:CG2	1:A:297:GLN:HG2	2.43	0.49
1:A:97:ASP:O	1:A:101:ASN:CG	2.56	0.49
1:A:62:VAL:HG22	1:A:108:HIS:CD2	2.47	0.49
1:A:151:VAL:HG22	1:A:152:LYS:N	2.28	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:76:ILE:C	1:A:176:VAL:HG23	2.38	0.48
1:A:129:VAL:C	1:A:130:THR:CA	2.79	0.48
1:A:64:ALA:O	1:A:65:LYS:C	2.57	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:98:ASN:O	1:A:99:ASP:C	2.57	0.48
1:A:50:LEU:CB	1:A:55:VAL:HG21	2.39	0.47
1:A:154:THR:C	1:A:156:ASP:N	2.70	0.47
1:A:111:VAL:CG1	1:A:112:ASP:N	2.77	0.47
1:A:163:TYR:CA	1:A:171:PHE:HD1	2.27	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
1:A:130:THR:HG1	1:A:138:ALA:C	2.22	0.46
2:B:144:MET:O	2:B:144:MET:HG2	2.14	0.46
1:A:117:LYS:O	1:A:120:LEU:HB3	2.15	0.46
1:A:240:LEU:O	1:A:243:LYS:HB2	2.14	0.46
1:A:210:THR:OG1	1:A:213:ASP:O	2.26	0.46
1:A:151:VAL:HG23	1:A:160:ALA:C	2.40	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:HIS:CD2	1:A:299:GLY:O	2.68	0.46
1:A:20:PRO:HB3	1:A:52:ALA:HB2	1.98	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: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:65:LYS:HZ3	4:A:893:POP:P2	2.39	0.45
1:A:193:ALA:HA	1:A:300:THR:HG21	1.99	0.45
2:B:143:GLN:C	2:B:145:MET:N	2.74	0.45
2:B:93:ASP:OD1	2:B:97:ASN:O	2.35	0.45
1:A:178:GLY:HA2	1:A:184:PRO:HA	1.98	0.45
1:A:89:ALA:O	1:A:92:VAL:N	2.49	0.45
1:A:315:VAL:HG21	1:A:325:LEU:CD1	2.40	0.45
1:A:150:ARG:HG3	1:A:162:GLN:HB2	1.99	0.44
1:A:163:TYR:C	1:A:171:PHE:CD1	2.95	0.44
1:A:253:VAL:HG11	2:B:87:GLU:HB3	1.99	0.44
1:A:313:ILE:HD11	1:A:315:VAL:HG22	1.99	0.44
1:A:66:SER:O	1:A:84:LYS:HG2	2.17	0.44
1:A:88:ARG:HH21	1:A:92:VAL:HG21	1.83	0.44
1:A:179:ASN:HB3	1:A:183:ILE:N	2.32	0.44
1:A:72:GLN:HA	1:A:113:LEU:HD21	1.99	0.44
1:A:104:LEU:HD23	1:A:108:HIS:O	2.18	0.44
1:A:130:THR:O	1:A:131:GLY:CA	2.59	0.44
1:A:346:GLU:HG3	1:A:347:ASN:O	2.18	0.44
1:A:150:ARG:NH1	1:A:150:ARG:HG2	2.33	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:56:ALA:O	1:A:186:THR:HG22	2.18	0.43
1:A:85:LEU:CD2	1:A:88:ARG:HG3	2.48	0.43
1:A:98:ASN:CA	1:A:101:ASN:ND2	2.73	0.43
1:A:81:ASN:C	1:A:82:LEU:HD23	2.44	0.43
1:A:297:GLN:HE21	1:A:297:GLN:CA	2.09	0.43
1:A:71:LEU:CD1	1:A:161:VAL:HG13	2.48	0.43
1:A:191:MET:HB3	1:A:191:MET:HE2	1.73	0.43
1:A:269:ILE:O	1:A:269:ILE:CG2	2.66	0.43
1:A:334:ILE:O	1:A:338:ARG:HG3	2.18	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:85:LEU:HD22	1:A:92:VAL:HG11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:ARG:HG2	1:A:171:PHE:CE1	2.55	0.42
1:A:62:VAL:O	1:A:63:HIS:ND1	2.53	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
1:A:18:GLY:N	6:A:925:HOH:O	2.52	0.41
1:A:135:GLY:O	1:A:137:VAL:HG13	2.19	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:11:ASN:O	1:A:15:ARG:HD2	2.20	0.41
1:A:39:MET:HB3	1:A:313:ILE:CD1	2.51	0.41
1:A:88:ARG:HE	1:A:88:ARG:HB3	1.66	0.40
1:A:288:HIS:HA	1:A:292:ALA:O	2.21	0.40
2:B:106:ARG:CZ	2:B:106:ARG:HB3	2.50	0.40
1:A:193:ALA:HB2	1:A:302:GLN:HG2	2.03	0.40
1:A:246:ARG:HD2	1:A:246:ARG:HA	1.74	0.40
1:A:300:THR:CG2	1:A:302:GLN:HB2	2.51	0.40
2:B:109:MET:HE3	2:B:109:MET:HA	2.03	0.40
1:A:11:ASN:HD21	1:A:14:ASP:H	1.67	0.40
1:A:103:SER:O	1:A:108:HIS:CB	2.55	0.40
1:A:130:THR:OG1	1:A:138:ALA:HB3	2.19	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 ⓘ

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	345/356 (97%)	301 (87%)	34 (10%)	10 (3%)	3 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	65/67 (97%)	60 (92%)	3 (5%)	2 (3%)	3	1
All	All	410/423 (97%)	361 (88%)	37 (9%)	12 (3%)	3	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%)	241 (89%)	30 (11%)	6	6
2	B	58/58 (100%)	51 (88%)	7 (12%)	5	4
All	All	329/332 (99%)	292 (89%)	37 (11%)	6	5

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	32	LYS
1	A	37	THR
1	A	57	THR

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Mol	Chain	Res	Type
1	A	63	HIS
1	A	82	LEU
1	A	91	GLU
1	A	119	ARG
1	A	123	LEU
1	A	136	VAL
1	A	137	VAL
1	A	146	GLN
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	246	ARG
1	A	256	GLU
1	A	285	ARG
1	A	297	GLN
1	A	300	THR
1	A	313	ILE
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	142	VAL
2	B	143	GLN

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	11	ASN
1	A	35	ASN
1	A	98	ASN
1	A	101	ASN
1	A	162	GLN

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Mol	Chain	Res	Type
1	A	197	HIS
1	A	260	GLN
1	A	293	GLN
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	135	GLN
2	B	137	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 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	$\# Z > 2$
4	POP	A	893	-	6,8,8	0.69	0	12,13,13	1.43	3 (25%)

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
4	POP	A	893	-	-	3/6/6/6	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	893	POP	O5-P2-O4	-2.70	100.32	110.83
4	A	893	POP	O6-P2-O4	2.23	119.51	110.83
4	A	893	POP	O6-P2-O	2.10	111.68	104.64

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	893	POP	P2-O-P1-O3
4	A	893	POP	P2-O-P1-O2
4	A	893	POP	P2-O-P1-O1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	893	POP	2	0

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	5

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	129:VAL	C	130:THR	N	1.71
1	A	85:LEU	C	86:PHE	N	1.63
1	A	133:ALA	C	134:ASP	N	0.92
1	A	86:PHE	C	87:GLY	N	0.78
1	A	130:THR	C	131:GLY	N	0.71

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%)	0.96	48 (13%) 6 5	7, 26, 47, 56	0
2	B	67/67 (100%)	0.64	2 (2%) 52 49	11, 22, 35, 47	0
All	All	416/423 (98%)	0.91	50 (12%) 9 6	7, 25, 47, 56	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	GLY	7.5
1	A	133	ALA	4.8
1	A	130	THR	4.5
1	A	173	ALA	4.5
1	A	174	VAL	3.7
1	A	115	LEU	3.7
1	A	78	VAL	3.5
1	A	132	MET	3.3
1	A	134	ASP	3.3
1	A	362	LEU	3.3
1	A	176	VAL	3.2
1	A	140	ASN	3.2
1	A	114	THR	3.2
1	A	7	ALA	3.1
1	A	361	GLY	3.1
1	A	63	HIS	3.1
1	A	165	ARG	3.0
1	A	171	PHE	3.0
2	B	145	MET	2.8
1	A	160	ALA	2.8
1	A	86	PHE	2.8
1	A	137	VAL	2.7
1	A	168	GLY	2.7
1	A	101	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	151	VAL	2.6
2	B	101	SER	2.5
1	A	98	ASN	2.5
1	A	87	GLY	2.5
1	A	135	GLY	2.5
1	A	85	LEU	2.5
1	A	180	ALA	2.5
1	A	54	GLY	2.5
1	A	136	VAL	2.4
1	A	170	ASP	2.4
1	A	110	ALA	2.4
1	A	149	PHE	2.4
1	A	225	ALA	2.4
1	A	89	ALA	2.3
1	A	310	ASP	2.3
1	A	159	TYR	2.3
1	A	129	VAL	2.3
1	A	107	GLY	2.2
1	A	220	ALA	2.2
1	A	183	ILE	2.1
1	A	106	HIS	2.1
1	A	150	ARG	2.1
1	A	142	ALA	2.1
1	A	108	HIS	2.0
1	A	113	LEU	2.0
1	A	357	LEU	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
4	POP	A	893	9/9	0.48	0.17	71,72,72,73	0
3	MG	A	908	1/1	0.82	0.12	23,23,23,23	0
5	CA	B	801	1/1	0.87	0.09	33,33,33,33	0
3	MG	A	907	1/1	0.90	0.17	21,21,21,21	0
5	CA	B	800	1/1	0.96	0.04	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.