



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 12:44 am BST

PDB ID : 1EX6
Title : CRYSTAL STRUCTURE OF UNLIGANDED FORM OF GUANYLATE KINASE FROM YEAST
Authors : Blaszczyk, J.; Ji, X.
Deposited on : 2000-05-01
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

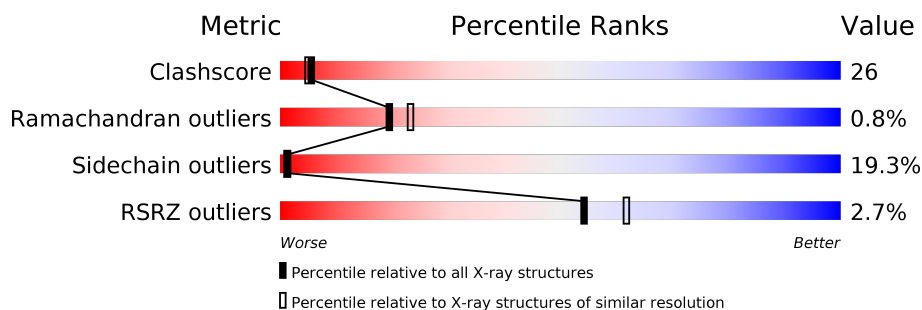
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	186	<div> <div></div> <div>52% 40% 8%</div> </div>
1	B	186	<div> <div>4%</div> <div>46% 41% 13%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3316 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 GUANYLATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	0	0	0
			1446	916	241	286	3			
1	B	186	Total	C	N	O	S	0	0	0
			1446	916	241	286	3			

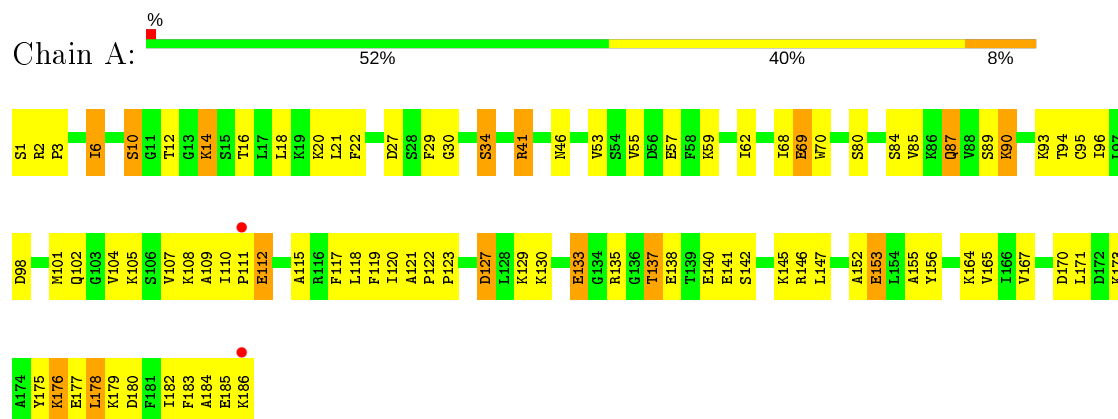
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	217	Total	O	0	0
			217	217		
2	B	207	Total	O	0	0
			207	207		

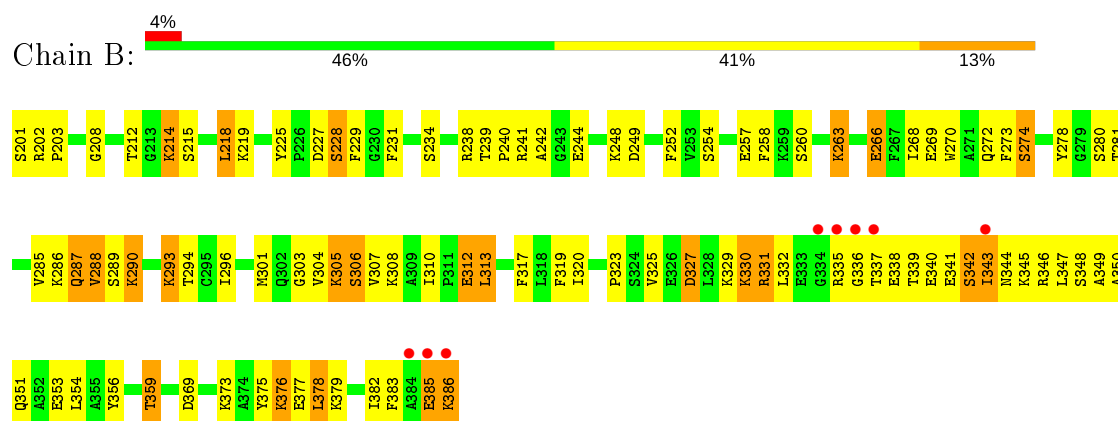
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: GUANYLATE KINASE



• Molecule 1: GUANYLATE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	51.56Å 51.56Å 119.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.53 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.7 (20.00-2.30) 81.8 (19.53-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.30Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.175 , 0.259 0.168 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 141.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l 0.093 for h,-h-k,-l 0.044 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3316	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.32	0/1472	0.85	0/1980
1	B	0.31	0/1472	0.85	2/1980 (0.1%)
All	All	0.32	0/2944	0.85	2/3960 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	241	ARG	N-CA-CB	5.26	120.08	110.60
1	B	331	ARG	NE-CZ-NH1	-5.19	117.70	120.30

There are no chirality outliers.

There are no planarity outliers.

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	1446	0	1445	68	0
1	B	1446	0	1442	82	0
2	A	217	0	0	9	0
2	B	207	0	0	6	0
All	All	3316	0	2887	149	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 26.

All (149) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:VAL:HG13	1:B:313:LEU:HD13	1.48	0.95
1:B:339:THR:HG22	1:B:342:SER:HB2	1.46	0.95
1:B:337:THR:HG22	1:B:337:THR:O	1.81	0.80
1:B:349:ALA:O	1:B:353:GLU:HG3	1.83	0.79
1:A:178:LEU:O	1:A:182:ILE:HG13	1.84	0.78
1:B:201:SER:HB3	1:B:385:GLU:OE2	1.85	0.77
1:A:127:ASP:HA	1:A:130:LYS:HG2	1.65	0.77
1:B:260:SER:HA	1:B:263:LYS:HG3	1.65	0.76
1:B:341:GLU:O	1:B:345:LYS:HD2	1.87	0.75
1:B:327:ASP:O	1:B:331:ARG:HG3	1.84	0.75
1:B:327:ASP:O	1:B:330:LYS:HG3	1.87	0.75
1:B:304:VAL:O	1:B:308:LYS:HG3	1.86	0.74
1:B:356:TYR:O	1:B:359:THR:HG23	1.87	0.74
1:A:55:VAL:O	1:A:59:LYS:HG3	1.88	0.73
1:B:373:LYS:O	1:B:376:LYS:HG3	1.89	0.73
1:B:335:ARG:HH22	1:B:346:ARG:NH2	1.86	0.72
1:B:342:SER:O	1:B:346:ARG:HG3	1.89	0.71
1:A:129:LYS:O	1:A:133:GLU:HG2	1.91	0.71
1:B:343:ILE:O	1:B:347:LEU:HG	1.90	0.71
1:B:266:GLU:HG2	2:B:523:HOH:O	1.90	0.70
1:A:140:GLU:HG3	1:A:141:GLU:H	1.57	0.70
1:B:332:LEU:HD22	1:B:346:ARG:HD3	1.73	0.70
1:B:254:SER:O	1:B:257:GLU:HB2	1.93	0.69
1:A:14:LYS:HE3	1:A:98:ASP:OD1	1.94	0.68
1:B:376:LYS:HE3	1:B:377:GLU:HG3	1.76	0.67
1:B:254:SER:OG	1:B:257:GLU:HG3	1.95	0.67
1:A:129:LYS:HE2	2:A:583:HOH:O	1.95	0.66
1:B:338:GLU:OE2	1:B:342:SER:HB3	1.95	0.66
1:B:225:TYR:HB3	1:B:228:SER:HB2	1.78	0.65
1:A:129:LYS:HB2	1:A:147:LEU:HD11	1.78	0.65
1:A:16:THR:O	1:A:20:LYS:HG3	1.97	0.64
1:B:339:THR:O	1:B:343:ILE:HB	1.98	0.64
1:B:281:THR:HB	2:B:523:HOH:O	1.98	0.64
1:A:34:SER:HB2	2:A:616:HOH:O	1.98	0.63
1:B:287:GLN:HG2	1:B:290:LYS:HZ3	1.64	0.62
1:B:338:GLU:HG3	1:B:339:THR:H	1.64	0.61
1:A:6:ILE:HG13	1:A:118:LEU:HD23	1.81	0.61
1:B:378:LEU:O	1:B:382:ILE:HD12	2.01	0.61
1:B:218:LEU:HD12	2:B:533:HOH:O	2.00	0.60
1:B:202:ARG:HD3	1:B:386:LYS:OXT	2.01	0.60
1:A:104:VAL:O	1:A:108:LYS:HG3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:THR:CG2	1:B:342:SER:H	2.15	0.60
1:A:12:THR:HG22	1:A:122:PRO:HB3	1.84	0.59
1:A:185:GLU:HG3	1:A:185:GLU:O	2.01	0.59
1:B:229:PHE:HD1	1:B:294:THR:HB	1.67	0.59
1:B:289:SER:HA	2:B:501:HOH:O	2.04	0.58
1:A:129:LYS:HB2	1:A:147:LEU:CD1	2.33	0.58
1:A:14:LYS:HG3	1:A:120:ILE:HG13	1.86	0.58
1:A:21:LEU:HD22	1:A:178:LEU:HD21	1.85	0.58
1:A:10:SER:HB2	2:A:428:HOH:O	2.04	0.57
1:B:301:MET:O	1:B:305:LYS:HD2	2.04	0.57
1:A:87:GLN:HG3	1:A:90:LYS:CE	2.33	0.57
1:B:350:ALA:O	1:B:354:LEU:HG	2.04	0.57
1:B:339:THR:HG22	1:B:342:SER:H	1.69	0.56
1:B:337:THR:CG2	1:B:337:THR:O	2.53	0.55
1:A:30:GLY:O	1:A:95:CYS:HA	2.07	0.55
1:A:3:PRO:HG2	1:A:115:ALA:HA	1.89	0.55
1:B:214:LYS:HG3	1:B:320:ILE:HG13	1.88	0.55
1:B:310:ILE:HG22	1:B:312:GLU:HG2	1.89	0.55
1:A:104:VAL:HG13	1:A:117:PHE:CE2	2.43	0.54
1:A:127:ASP:OD1	1:A:130:LYS:HE3	2.07	0.54
1:B:338:GLU:HG2	1:B:343:ILE:HG13	1.90	0.54
1:A:14:LYS:HG3	1:A:120:ILE:CG1	2.38	0.54
1:A:105:LYS:HE3	1:A:156:TYR:CE1	2.42	0.53
1:B:229:PHE:CD1	1:B:294:THR:HB	2.42	0.53
1:B:268:ILE:HG12	1:B:313:LEU:HD11	1.90	0.53
1:B:240:PRO:HG3	1:B:252:PHE:CD2	2.44	0.52
1:A:177:GLU:HG2	2:A:479:HOH:O	2.09	0.52
1:A:127:ASP:O	1:A:130:LYS:HG2	2.11	0.51
1:A:129:LYS:HA	1:A:147:LEU:HD21	1.93	0.51
1:B:260:SER:HA	1:B:263:LYS:CG	2.39	0.50
1:A:185:GLU:HB3	2:A:401:HOH:O	2.10	0.50
1:A:127:ASP:CA	1:A:130:LYS:HG2	2.37	0.50
1:A:69:GLU:OE2	1:A:80:SER:OG	2.26	0.50
1:B:375:TYR:CZ	1:B:379:LYS:HD3	2.48	0.49
1:A:22:PHE:CE1	1:A:96:ILE:HG13	2.47	0.49
1:A:104:VAL:HG13	1:A:117:PHE:CD2	2.48	0.49
1:B:287:GLN:O	1:B:290:LYS:HG3	2.12	0.49
1:B:307:VAL:HA	1:B:310:ILE:HD12	1.94	0.48
1:A:87:GLN:HG3	1:A:90:LYS:NZ	2.28	0.48
1:A:137:THR:O	1:A:137:THR:OG1	2.29	0.47
1:B:306:SER:O	1:B:310:ILE:HD12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ILE:HD11	1:A:85:VAL:HG21	1.96	0.47
1:B:238:ARG:HH11	1:B:238:ARG:HG2	1.80	0.47
1:B:286:LYS:O	1:B:290:LYS:HG2	2.14	0.47
1:B:375:TYR:O	1:B:379:LYS:HG3	2.14	0.47
1:B:335:ARG:HH22	1:B:346:ARG:HH21	1.61	0.47
1:B:356:TYR:C	1:B:359:THR:HG23	2.35	0.47
1:B:386:LYS:HE2	1:B:386:LYS:HB3	1.48	0.47
1:A:118:LEU:HD21	1:A:178:LEU:HD12	1.97	0.47
1:A:6:ILE:HG23	1:A:6:ILE:O	2.15	0.47
1:B:208:GLY:HA2	1:B:319:PHE:CE1	2.49	0.47
1:A:27:ASP:O	1:A:93:LYS:HG2	2.14	0.47
1:A:176:LYS:HB3	1:A:176:LYS:HE3	1.70	0.46
1:A:59:LYS:HE3	1:A:59:LYS:HB3	1.31	0.46
1:A:182:ILE:HG22	1:A:183:PHE:CD1	2.51	0.46
1:A:87:GLN:HG3	1:A:90:LYS:HE3	1.98	0.45
1:B:238:ARG:NH2	1:B:244:GLU:OE2	2.50	0.45
1:A:110:ILE:HG22	1:A:112:GLU:HG2	1.98	0.45
1:A:135:ARG:NH1	2:A:428:HOH:O	2.50	0.45
1:A:119:PHE:HD2	1:A:165:VAL:HG13	1.82	0.45
1:B:214:LYS:HB2	1:B:214:LYS:HE3	1.74	0.45
1:A:173:LYS:HE3	2:A:600:HOH:O	2.17	0.45
1:B:385:GLU:OE1	1:B:386:LYS:N	2.50	0.45
1:B:201:SER:N	2:B:501:HOH:O	2.50	0.45
1:B:335:ARG:NH2	1:B:338:GLU:OE1	2.50	0.45
1:A:130:LYS:HB3	2:A:749:HOH:O	2.16	0.44
1:B:238:ARG:NH1	1:B:239:THR:O	2.50	0.44
1:A:176:LYS:NZ	2:A:538:HOH:O	2.50	0.44
1:A:1:SER:OG	1:A:2:ARG:N	2.49	0.44
1:B:231:PHE:HD1	1:B:296:ILE:O	2.00	0.44
1:A:18:LEU:CD2	1:A:96:ILE:HG21	2.48	0.44
1:A:171:LEU:O	1:A:175:TYR:HB2	2.18	0.44
1:A:141:GLU:OE1	1:A:145:LYS:NZ	2.50	0.44
1:B:270:TRP:HA	1:B:278:TYR:O	2.18	0.44
1:B:339:THR:HG23	1:B:340:GLU:N	2.33	0.44
1:A:109:ALA:C	1:A:111:PRO:HD3	2.38	0.43
1:A:70:TRP:CH2	1:B:312:GLU:HB3	2.53	0.43
1:B:285:VAL:HG11	1:B:313:LEU:CD2	2.48	0.43
1:B:287:GLN:HA	1:B:290:LYS:HG2	2.00	0.43
1:B:308:LYS:HG2	1:B:317:PHE:CZ	2.53	0.43
1:B:202:ARG:HH11	1:B:386:LYS:C	2.22	0.43
1:A:133:GLU:HG2	1:A:133:GLU:H	1.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:SER:HB2	2:B:632:HOH:O	2.18	0.43
1:B:303:GLY:O	1:B:307:VAL:HG23	2.18	0.43
1:A:22:PHE:HE1	1:A:96:ILE:HG13	1.84	0.42
1:B:332:LEU:HD22	1:B:346:ARG:CD	2.45	0.42
1:B:248:LYS:HG2	1:B:249:ASP:N	2.35	0.42
1:B:273:PHE:HD2	1:B:278:TYR:CD1	2.37	0.42
1:A:119:PHE:CE2	1:A:121:ALA:HB2	2.55	0.41
1:A:53:VAL:HB	1:A:57:GLU:HB2	2.02	0.41
1:A:180:ASP:O	1:A:184:ALA:HB2	2.20	0.41
1:B:373:LYS:O	1:B:376:LYS:HE3	2.20	0.41
1:B:382:ILE:HG22	1:B:383:PHE:CD1	2.55	0.41
1:A:101:MET:HB2	1:A:153:GLU:HG3	2.03	0.41
1:A:41:ARG:HD2	1:A:41:ARG:HA	1.27	0.41
1:A:123:PRO:HA	1:A:167:VAL:CG1	2.50	0.41
1:A:152:ALA:O	1:A:155:ALA:HB3	2.21	0.41
1:A:90:LYS:HE3	1:A:90:LYS:HB3	1.78	0.41
1:B:285:VAL:O	1:B:288:VAL:HG23	2.20	0.41
1:B:323:PRO:HG3	1:B:369:ASP:HA	2.03	0.41
1:B:327:ASP:HA	1:B:330:LYS:CG	2.50	0.41
1:A:62:ILE:HD13	1:A:70:TRP:CZ2	2.56	0.41
1:B:202:ARG:HA	1:B:203:PRO:HD3	1.90	0.41
1:A:107:VAL:HG12	1:A:117:PHE:HZ	1.86	0.40
1:B:227:ASP:O	1:B:293:LYS:HD2	2.20	0.40
1:A:29:PHE:CD1	1:A:96:ILE:HD11	2.57	0.40
1:B:268:ILE:HB	1:B:280:SER:O	2.21	0.40
1:B:330:LYS:HE3	1:B:330:LYS:HB2	1.34	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	184/186 (99%)	174 (95%)	9 (5%)	1 (0%)	29	35
1	B	184/186 (99%)	174 (95%)	8 (4%)	2 (1%)	14	15
All	All	368/372 (99%)	348 (95%)	17 (5%)	3 (1%)	19	23

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	242	ALA
1	B	336	GLY
1	A	170	ASP

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	158/158 (100%)	132 (84%)	26 (16%)	2	2
1	B	158/158 (100%)	123 (78%)	35 (22%)	1	1
All	All	316/316 (100%)	255 (81%)	61 (19%)	1	1

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ILE
1	A	10	SER
1	A	14	LYS
1	A	34	SER
1	A	41	ARG
1	A	46	ASN
1	A	69	GLU
1	A	84	SER
1	A	87	GLN
1	A	89	SER
1	A	90	LYS
1	A	94	THR
1	A	102	GLN

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Mol	Chain	Res	Type
1	A	112	GLU
1	A	127	ASP
1	A	133	GLU
1	A	137	THR
1	A	138	GLU
1	A	142	SER
1	A	146	ARG
1	A	153	GLU
1	A	164	LYS
1	A	176	LYS
1	A	178	LEU
1	A	179	LYS
1	A	186	LYS
1	B	212	THR
1	B	214	LYS
1	B	215	SER
1	B	218	LEU
1	B	219	LYS
1	B	228	SER
1	B	234	SER
1	B	258	PHE
1	B	263	LYS
1	B	266	GLU
1	B	269	GLU
1	B	272	GLN
1	B	274	SER
1	B	287	GLN
1	B	288	VAL
1	B	290	LYS
1	B	293	LYS
1	B	305	LYS
1	B	306	SER
1	B	312	GLU
1	B	313	LEU
1	B	325	VAL
1	B	327	ASP
1	B	329	LYS
1	B	330	LYS
1	B	342	SER
1	B	343	ILE
1	B	344	ASN
1	B	348	SER

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Mol	Chain	Res	Type
1	B	351	GLN
1	B	359	THR
1	B	376	LYS
1	B	378	LEU
1	B	385	GLU
1	B	386	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	264	ASN
1	B	265	ASN
1	B	272	GLN
1	B	287	GLN
1	B	344	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	186/186 (100%)	-0.46	2 (1%) 80 85	10, 29, 45, 69	0
1	B	186/186 (100%)	-0.17	8 (4%) 35 42	12, 29, 56, 68	3 (1%)
All	All	372/372 (100%)	-0.31	10 (2%) 54 62	10, 29, 53, 69	3 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	337	THR	5.7
1	B	336	GLY	5.7
1	B	386	LYS	3.4
1	A	111	PRO	3.1
1	B	335	ARG	3.1
1	B	385	GLU	2.8
1	B	343	ILE	2.8
1	B	334	GLY	2.7
1	A	186	LYS	2.6
1	B	384	ALA	2.5

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](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.