



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:00 am BST

PDB ID : 4JZ7
Title : Carbamate kinase from Giardia lamblia bound to AMP-PNP
Authors : Lim, K.; Herzberg, O.
Deposited on : 2013-04-02
Resolution : 2.60 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
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

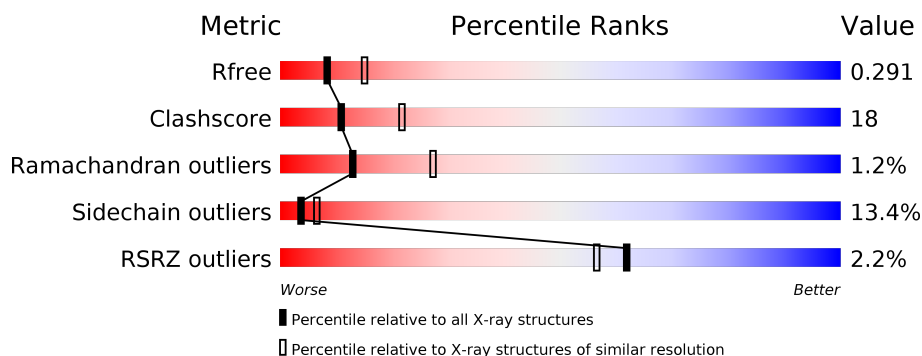
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	317	
1	B	317	
1	C	317	
1	D	317	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9567 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 Carbamate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2118	1329	360	411	18			
1	B	285	Total	C	N	O	S	0	0	0
			2118	1329	360	411	18			
1	C	316	Total	C	N	O	S	0	0	0
			2366	1483	408	456	19			
1	D	316	Total	C	N	O	S	0	0	0
			2366	1483	408	456	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	INSERTION	UNP A8BB85
B	0	GLY	-	INSERTION	UNP A8BB85
C	0	GLY	-	INSERTION	UNP A8BB85
D	0	GLY	-	INSERTION	UNP A8BB85

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	B	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	C	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	D	1	Total 31	C 10	N 6	O 12	P 3	0	0

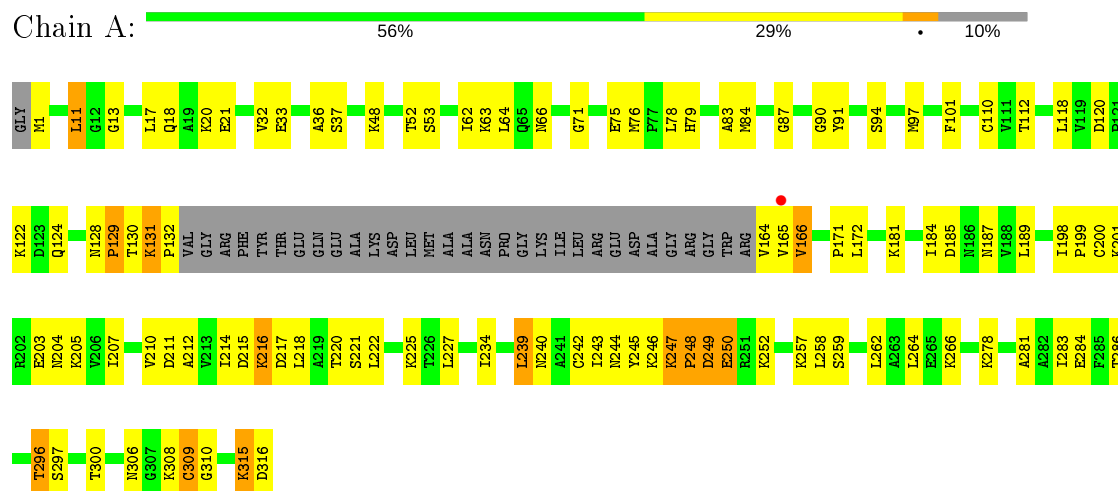
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	128	Total O 128 128	0	0
3	B	101	Total O 101 101	0	0
3	C	119	Total O 119 119	0	0
3	D	127	Total O 127 127	0	0

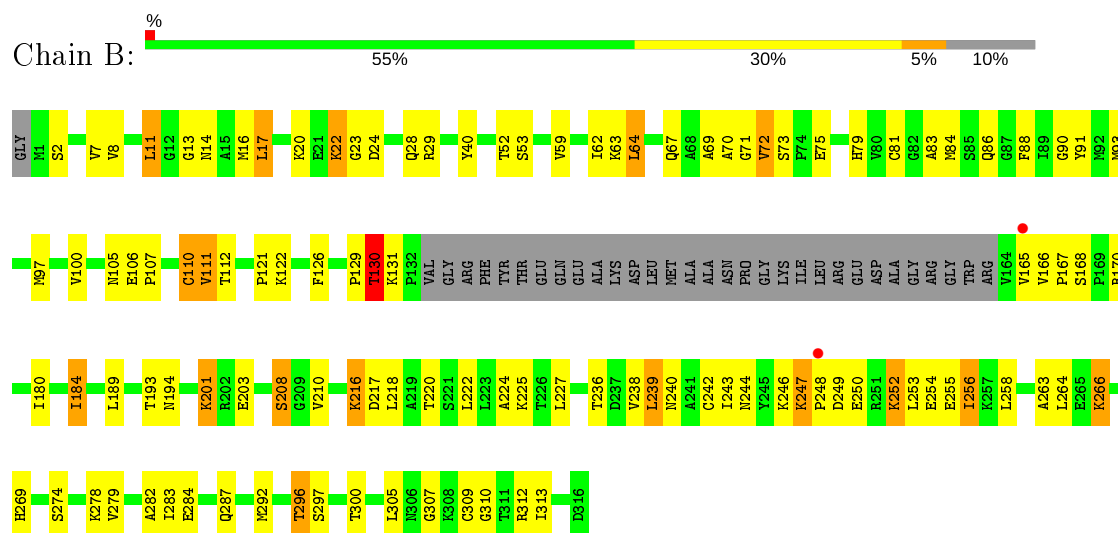
3 Residue-property plots

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: Carbamate kinase

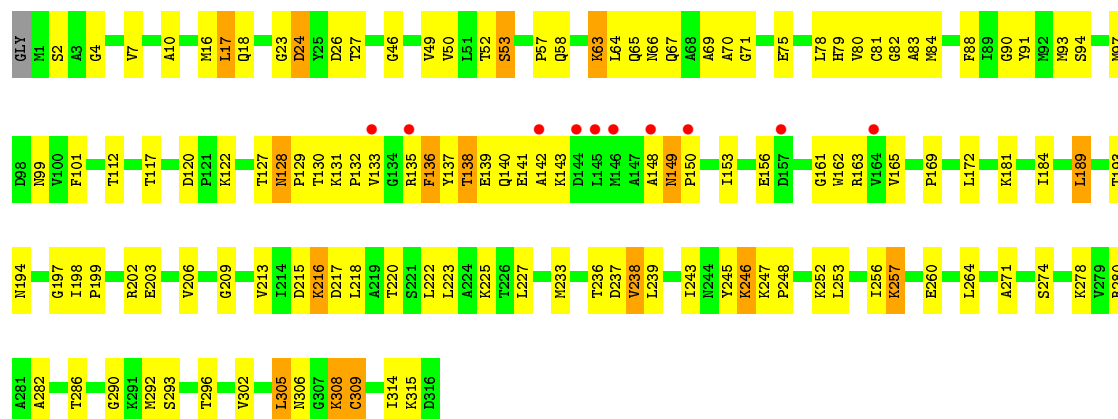


• Molecule 1: Carbamate kinase

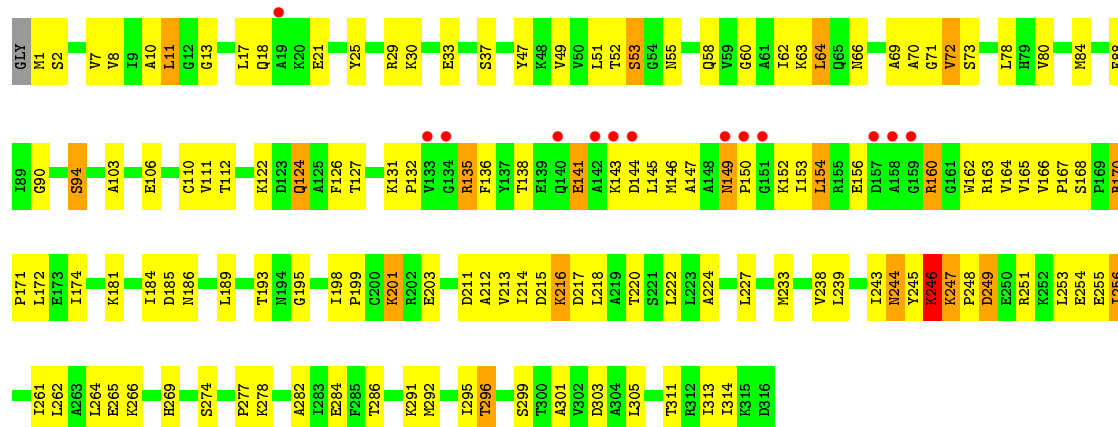


• Molecule 1: Carbamate kinase





• Molecule 1: Carbamate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.63 Å 97.05 Å 102.12 Å 90.00° 106.69° 90.00°	Depositor
Resolution (Å)	48.90 – 2.60 48.91 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (48.90-2.60) 99.3 (48.91-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.66 (at 2.61 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.217 , 0.292 0.211 , 0.291	Depositor DCC
R_{free} test set	1616 reflections (4.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.1	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9567	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.5655e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ANP

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	0.49	0/2144	0.66	0/2898
1	B	0.47	0/2144	0.63	0/2898
1	C	0.53	0/2398	0.63	0/3240
1	D	0.49	0/2398	0.66	0/3240
All	All	0.50	0/9084	0.65	0/12276

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	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	315	LYS	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	2118	0	2186	72	0
1	B	2118	0	2186	80	0
1	C	2366	0	2430	88	0
1	D	2366	0	2430	105	0
2	A	31	0	13	1	0
2	B	31	0	13	6	0
2	C	31	0	13	5	0
2	D	31	0	13	4	0
3	A	128	0	0	5	0
3	B	101	0	0	9	0
3	C	119	0	0	6	0
3	D	127	0	0	12	0
All	All	9567	0	9284	335	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 18.

All (335) 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:C:216:LYS:HE2	1:C:217:ASP:OD1	1.59	1.01
1:B:67:GLN:HG3	1:B:75:GLU:OE1	1.71	0.91
1:A:243:ILE:HG21	1:A:264:LEU:HD13	1.57	0.87
1:B:131:LYS:O	1:B:166:VAL:HG22	1.77	0.84
1:A:315:LYS:HB2	3:A:1112:HOH:O	1.78	0.83
1:B:105:ASN:HB3	3:B:592:HOH:O	1.80	0.81
1:B:52:THR:HG21	1:B:220:THR:OG1	1.82	0.80
1:A:164:VAL:HG12	1:A:165:VAL:H	1.47	0.78
1:A:225:LYS:HE2	3:A:1220:HOH:O	1.83	0.77
1:B:131:LYS:N	1:B:166:VAL:O	2.17	0.77
1:A:52:THR:HG21	1:A:220:THR:OG1	1.85	0.77
1:B:90:GLY:HA3	1:B:112:THR:HG21	1.67	0.77
1:D:52:THR:HG21	1:D:220:THR:OG1	1.85	0.76
1:C:136:PHE:HB3	1:C:161:GLY:HA3	1.67	0.76
1:B:258:LEU:HD21	1:B:287:GLN:HG2	1.70	0.73
1:C:52:THR:HG21	1:C:220:THR:OG1	1.89	0.72
1:C:127:THR:O	1:C:129:PRO:HD3	1.90	0.72
1:B:255:GLU:C	1:B:256:ILE:HD13	2.10	0.71
1:D:90:GLY:HA3	1:D:112:THR:HG21	1.72	0.71
1:D:256:ILE:HD11	1:D:311:THR:HG23	1.72	0.71
1:D:181:LYS:NZ	1:D:185:ASP:OD2	2.24	0.70
1:B:67:GLN:HB3	1:D:135:ARG:NH2	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:VAL:HB	1:C:189:LEU:HD13	1.75	0.69
1:C:90:GLY:HA3	1:C:112:THR:HG21	1.76	0.68
1:C:99:ASN:HA	3:C:509:HOH:O	1.94	0.68
1:D:154:LEU:HD12	1:D:162:TRP:HB3	1.75	0.67
1:C:156:GLU:HB2	1:C:162:TRP:CE2	2.30	0.67
1:D:216:LYS:HE2	1:D:217:ASP:OD1	1.95	0.67
1:B:216:LYS:HE2	1:B:217:ASP:OD1	1.95	0.66
1:A:129:PRO:CG	1:A:165:VAL:HG13	2.27	0.65
1:A:90:GLY:HA3	1:A:112:THR:HG21	1.79	0.65
1:B:72:VAL:O	1:D:160:ARG:NH2	2.25	0.64
1:C:236:THR:HG23	1:C:238:VAL:H	1.62	0.64
1:D:199:PRO:HG3	1:D:214:ILE:HG13	1.80	0.64
1:B:130:THR:N	1:B:166:VAL:O	2.29	0.64
1:C:138:THR:HG22	1:C:140:GLN:H	1.63	0.64
1:C:82:GLY:HA3	3:C:525:HOH:O	1.98	0.64
1:D:131:LYS:HE3	3:D:615:HOH:O	1.99	0.64
1:B:249:ASP:N	1:B:249:ASP:OD2	2.31	0.63
1:A:129:PRO:HG2	1:A:165:VAL:HG13	1.80	0.63
1:D:249:ASP:OD2	1:D:249:ASP:N	2.31	0.63
1:C:93:MET:O	1:C:97:MET:HB2	1.99	0.62
1:A:66:ASN:O	1:A:75:GLU:HG3	1.99	0.62
1:B:243:ILE:HG22	1:B:269:HIS:CG	2.34	0.62
1:B:222:LEU:HA	1:B:225:LYS:HE2	1.82	0.62
1:A:245:TYR:CE2	1:A:246:LYS:HG2	2.35	0.62
1:A:234:ILE:HG21	1:A:278:LYS:HD3	1.81	0.61
1:C:49:VAL:HB	1:C:189:LEU:CD1	2.30	0.61
1:A:308:LYS:O	1:A:309:CYS:HB3	1.99	0.61
2:B:401:ANP:N3	2:B:401:ANP:H2'	2.14	0.61
1:C:84:MET:HB3	1:D:84:MET:HG2	1.83	0.61
1:D:55:ASN:ND2	1:D:195:GLY:HA3	2.16	0.61
1:C:215:ASP:HB3	1:C:218:LEU:HD12	1.82	0.61
1:B:22:LYS:O	1:B:24:ASP:N	2.30	0.60
1:B:75:GLU:OE2	1:D:141:GLU:OE2	2.19	0.60
1:D:292:MET:CE	1:D:314:ILE:HD11	2.32	0.60
1:D:69:ALA:O	1:D:71:GLY:N	2.34	0.60
1:D:201:LYS:HD3	3:D:518:HOH:O	2.02	0.59
1:B:236:THR:HG23	1:B:238:VAL:H	1.67	0.59
1:A:242:CYS:SG	1:A:252:LYS:HD3	2.42	0.59
1:A:79:HIS:HB3	1:A:210:VAL:O	2.01	0.59
1:C:247:LYS:HB3	1:C:248:PRO:HD2	1.84	0.59
1:D:261:ILE:HA	1:D:264:LEU:HB2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:ILE:CG2	1:A:264:LEU:HD13	2.31	0.59
1:B:296:THR:HG23	1:B:310:GLY:HA3	1.84	0.59
1:C:278:LYS:NZ	2:C:401:ANP:O2B	2.33	0.59
1:B:40:TYR:OH	1:B:106:GLU:OE2	2.17	0.59
1:B:121:PRO:O	1:B:126:PHE:CD2	2.55	0.58
1:C:148:ALA:O	1:C:150:PRO:HD3	2.03	0.58
1:C:67:GLN:HG2	1:C:75:GLU:OE1	2.03	0.58
1:C:314:ILE:HG22	1:C:315:LYS:H	1.68	0.58
1:B:278:LYS:NZ	2:B:401:ANP:O2B	2.32	0.58
1:D:244:ASN:ND2	1:D:249:ASP:O	2.36	0.58
1:D:69:ALA:C	1:D:71:GLY:H	2.06	0.58
1:B:263:ALA:O	1:B:266:LYS:HG3	2.03	0.58
1:C:292:MET:HG3	1:C:314:ILE:CD1	2.34	0.57
1:C:120:ASP:HB2	1:C:172:LEU:HD11	1.86	0.57
1:D:78:LEU:N	1:D:211:ASP:OD2	2.35	0.57
1:D:149:ASN:N	1:D:149:ASN:OD1	2.38	0.57
1:B:13:GLY:H	2:B:401:ANP:HNB1	1.51	0.56
1:A:306:ASN:O	1:A:308:LYS:HE3	2.05	0.56
1:D:153:ILE:O	1:D:164:VAL:HA	2.05	0.56
1:C:128:ASN:O	1:C:130:THR:HG23	2.04	0.56
1:D:243:ILE:HD11	1:D:251:ARG:HD3	1.87	0.56
1:D:282:ALA:HB1	1:D:313:ILE:HD12	1.88	0.56
1:B:282:ALA:HB1	1:B:313:ILE:HD12	1.87	0.56
1:C:292:MET:HG3	1:C:314:ILE:HD13	1.87	0.56
1:D:143:LYS:HA	1:D:146:MET:HE2	1.88	0.56
1:C:16:MET:O	1:C:17:LEU:HD13	2.07	0.55
1:D:295:ILE:O	1:D:296:THR:HG22	2.07	0.55
1:A:259:SER:HB3	1:A:316:ASP:OD2	2.07	0.55
1:B:258:LEU:HD21	1:B:287:GLN:CG	2.37	0.55
1:B:279:VAL:HG12	1:B:283:ILE:HD12	1.89	0.55
1:B:97:MET:O	1:B:100:VAL:HG12	2.07	0.55
1:D:244:ASN:HB3	1:D:247:LYS:HD2	1.89	0.55
1:C:80:VAL:HG13	1:D:88:PHE:CD1	2.42	0.55
1:D:49:VAL:HB	1:D:189:LEU:CD2	2.37	0.55
1:A:216:LYS:HE2	1:A:217:ASP:OD1	2.08	0.54
1:D:274:SER:O	1:D:277:PRO:HD2	2.08	0.54
1:A:239:LEU:HG	1:A:240:ASN:OD1	2.07	0.54
1:A:244:ASN:O	1:A:247:LYS:HB2	2.08	0.54
1:C:137:TYR:HB2	1:C:142:ALA:HB2	1.89	0.54
1:D:110:CYS:HB3	3:D:559:HOH:O	2.07	0.54
1:A:129:PRO:HB3	1:A:166:VAL:N	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:GLY:HA2	1:C:135:ARG:NH1	2.23	0.54
1:A:201:LYS:HD2	1:A:203:GLU:HG3	1.90	0.54
1:A:129:PRO:CB	1:A:165:VAL:HG13	2.37	0.54
1:B:170:ARG:NH1	1:B:284:GLU:OE1	2.41	0.54
1:D:154:LEU:HB3	1:D:162:TRP:HE3	1.73	0.54
1:B:246:LYS:N	1:B:250:GLU:OE1	2.33	0.53
1:C:181:LYS:O	1:C:184:ILE:HG13	2.07	0.53
1:A:164:VAL:HG12	1:A:165:VAL:N	2.22	0.53
1:C:52:THR:HG22	1:C:53:SER:N	2.22	0.53
1:C:133:VAL:HA	3:C:599:HOH:O	2.08	0.53
1:D:138:THR:HG22	1:D:141:GLU:CG	2.38	0.53
1:D:278:LYS:NZ	2:D:401:ANP:O2B	2.30	0.53
1:A:13:GLY:H	2:A:1001:ANP:HNB1	1.57	0.53
1:D:80:VAL:HG12	1:D:84:MET:CE	2.38	0.53
1:D:138:THR:HG22	1:D:141:GLU:HG2	1.89	0.53
1:D:238:VAL:HG21	2:D:401:ANP:H2	1.91	0.53
1:A:262:LEU:HD23	1:A:283:ILE:HD13	1.91	0.53
1:C:256:ILE:HG13	1:C:260:GLU:OE2	2.09	0.53
1:D:292:MET:HE2	1:D:314:ILE:HD11	1.91	0.53
1:D:256:ILE:CD1	1:D:311:THR:HG23	2.38	0.53
1:A:201:LYS:O	1:A:207:ILE:HA	2.09	0.53
1:A:216:LYS:NZ	3:A:1104:HOH:O	2.40	0.53
1:B:8:VAL:HG21	1:B:224:ALA:HA	1.90	0.53
1:D:29:ARG:HH21	1:D:30:LYS:HB3	1.74	0.53
1:B:239:LEU:HD12	1:B:297:SER:HB3	1.91	0.52
1:C:26:ASP:HB2	3:C:566:HOH:O	2.08	0.52
1:C:131:LYS:HD2	1:C:132:PRO:HD2	1.92	0.52
1:D:215:ASP:HB3	1:D:218:LEU:HD12	1.92	0.52
1:B:69:ALA:C	1:B:71:GLY:H	2.13	0.52
1:C:314:ILE:HG22	1:C:315:LYS:N	2.23	0.52
1:D:145:LEU:HD22	1:D:149:ASN:OD1	2.10	0.52
1:D:174:ILE:HD12	1:D:222:LEU:HD23	1.91	0.52
1:D:135:ARG:HG2	1:D:136:PHE:N	2.24	0.51
1:A:234:ILE:CG2	1:A:278:LYS:HD3	2.40	0.51
1:A:296:THR:HG23	1:A:310:GLY:HA3	1.92	0.51
1:C:52:THR:CG2	1:C:53:SER:N	2.74	0.51
1:A:131:LYS:HG3	1:A:132:PRO:HD2	1.91	0.51
1:B:14:ASN:HA	1:B:17:LEU:O	2.11	0.51
1:C:138:THR:CG2	1:C:139:GLU:N	2.73	0.51
1:A:300:THR:OG1	1:A:300:THR:O	2.24	0.51
1:C:237:ASP:OD1	2:C:401:ANP:O3'	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:LEU:HA	1:C:225:LYS:HZ1	1.75	0.50
1:C:65:GLN:NE2	1:D:66:ASN:OD1	2.45	0.50
1:D:170:ARG:NH1	1:D:284:GLU:OE1	2.45	0.50
1:A:215:ASP:HB3	1:A:218:LEU:HD12	1.94	0.50
1:B:263:ALA:HA	1:B:266:LYS:HE3	1.94	0.50
1:B:67:GLN:CG	1:B:75:GLU:OE1	2.53	0.50
1:C:216:LYS:NZ	2:C:401:ANP:O2G	2.44	0.50
1:A:296:THR:HG23	1:A:310:GLY:CA	2.42	0.50
1:B:201:LYS:NZ	1:B:208:SER:OG	2.28	0.50
1:B:239:LEU:HB3	1:B:240:ASN:OD1	2.12	0.50
1:C:128:ASN:N	1:C:128:ASN:OD1	2.45	0.50
1:C:282:ALA:HB1	1:C:293:SER:HB3	1.93	0.50
1:A:83:ALA:HA	1:A:198:ILE:HD11	1.94	0.49
1:D:245:TYR:C	1:D:247:LYS:H	2.14	0.49
1:C:83:ALA:HA	1:C:198:ILE:HD11	1.95	0.49
1:D:1:MET:HE1	1:D:47:TYR:O	2.13	0.49
1:D:53:SER:O	1:D:216:LYS:HG3	2.12	0.49
1:D:286:THR:HG23	1:D:291:LYS:O	2.13	0.49
1:A:181:LYS:HD3	1:A:185:ASP:OD2	2.13	0.49
1:C:194:ASN:ND2	3:C:571:HOH:O	2.34	0.49
1:B:274:SER:HB2	2:B:401:ANP:O4'	2.13	0.48
2:D:401:ANP:N7	3:D:555:HOH:O	2.34	0.48
1:B:52:THR:CG2	1:B:53:SER:N	2.75	0.48
1:D:52:THR:CG2	1:D:53:SER:N	2.77	0.48
1:B:110:CYS:SG	1:B:111:VAL:N	2.86	0.48
1:B:121:PRO:HA	1:B:126:PHE:CE2	2.48	0.48
1:D:253:LEU:HB3	1:D:256:ILE:HD11	1.96	0.48
1:A:166:VAL:HG23	1:A:212:ALA:HA	1.95	0.48
1:A:18:GLN:HB2	1:A:21:GLU:OE2	2.14	0.48
1:C:91:TYR:CE1	1:D:198:ILE:HG21	2.48	0.48
1:C:169:PRO:HD2	1:C:213:VAL:O	2.13	0.48
1:A:53:SER:O	1:A:216:LYS:HG3	2.13	0.48
1:D:64:LEU:HD22	1:D:64:LEU:O	2.14	0.48
1:B:16:MET:O	1:B:17:LEU:HD13	2.14	0.48
1:D:18:GLN:O	1:D:21:GLU:HB2	2.14	0.48
1:B:201:LYS:HE2	3:B:555:HOH:O	2.14	0.47
1:B:263:ALA:HA	1:B:266:LYS:HG2	1.95	0.47
1:C:222:LEU:HA	1:C:225:LYS:NZ	2.29	0.47
1:A:130:THR:HG22	1:A:131:LYS:N	2.28	0.47
1:A:78:LEU:HD23	1:A:212:ALA:N	2.29	0.47
1:B:180:ILE:O	1:B:184:ILE:HG12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:147:ALA:O	1:D:150:PRO:HG3	2.14	0.47
1:A:63:LYS:HA	1:A:63:LYS:HD2	1.63	0.47
1:C:24:ASP:OD2	1:C:27:THR:N	2.34	0.47
1:D:184:ILE:HD11	3:D:611:HOH:O	2.14	0.47
1:C:308:LYS:O	1:C:309:CYS:HB3	2.13	0.47
1:D:49:VAL:HB	1:D:189:LEU:HD21	1.95	0.47
1:C:243:ILE:HG21	1:C:264:LEU:HD13	1.96	0.47
1:D:58:GLN:O	1:D:62:ILE:HG12	2.15	0.47
1:C:129:PRO:O	1:C:165:VAL:HB	2.14	0.47
1:B:307:GLY:HA2	3:B:554:HOH:O	2.15	0.47
2:B:401:ANP:O1G	3:B:549:HOH:O	2.21	0.47
1:C:91:TYR:CD1	1:D:198:ILE:HD13	2.50	0.47
1:B:88:PHE:CD1	1:B:88:PHE:C	2.88	0.46
1:C:306:ASN:O	1:C:308:LYS:HD3	2.15	0.46
1:D:254:GLU:HA	1:D:254:GLU:OE1	2.14	0.46
1:C:88:PHE:CD1	1:D:80:VAL:HG13	2.50	0.46
1:B:107:PRO:HG2	3:B:590:HOH:O	2.16	0.46
1:C:117:THR:OG1	1:C:197:GLY:HA3	2.15	0.46
1:C:50:VAL:HG11	1:C:223:LEU:HD21	1.98	0.46
1:C:286:THR:O	1:C:290:GLY:N	2.44	0.46
1:D:90:GLY:O	1:D:94:SER:HB2	2.16	0.46
1:C:10:ALA:HB2	1:C:220:THR:HG21	1.98	0.46
1:C:149:ASN:OD1	1:C:149:ASN:N	2.48	0.46
1:B:86:GLN:HG2	1:B:194:ASN:HB2	1.97	0.46
1:A:120:ASP:HB2	1:A:172:LEU:HD11	1.98	0.46
1:B:69:ALA:O	1:B:71:GLY:N	2.49	0.46
1:D:126:PHE:CD1	1:D:167:PRO:HG2	2.50	0.46
1:D:244:ASN:HB3	1:D:247:LYS:CD	2.46	0.46
2:C:401:ANP:N3	2:C:401:ANP:H2'	2.30	0.46
1:D:51:LEU:HD23	1:D:51:LEU:HA	1.59	0.46
1:B:300:THR:HG22	3:B:563:HOH:O	2.15	0.46
1:A:118:LEU:HD23	1:A:200:CYS:HB2	1.97	0.46
1:A:278:LYS:O	1:A:281:ALA:HB3	2.15	0.46
1:B:69:ALA:HB1	1:B:72:VAL:HG23	1.97	0.46
1:A:249:ASP:OD2	1:A:249:ASP:N	2.49	0.45
1:D:58:GLN:HA	3:D:550:HOH:O	2.15	0.45
1:D:60:GLY:HA2	3:D:569:HOH:O	2.15	0.45
1:D:170:ARG:HA	1:D:171:PRO:HD3	1.79	0.45
1:B:309:CYS:O	1:B:312:ARG:NE	2.50	0.45
1:B:88:PHE:O	1:B:91:TYR:HB3	2.17	0.45
1:A:204:ASN:O	1:A:205:LYS:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:LEU:O	1:B:216:LYS:HE3	2.16	0.45
1:D:299:SER:HB3	3:D:521:HOH:O	2.17	0.45
1:A:239:LEU:HA	1:A:297:SER:H	1.81	0.45
1:C:131:LYS:HG3	1:C:133:VAL:HG23	1.99	0.45
1:C:156:GLU:HG2	1:C:161:GLY:O	2.16	0.45
1:D:245:TYR:O	1:D:247:LYS:N	2.50	0.45
1:C:225:LYS:NZ	3:C:504:HOH:O	2.48	0.45
1:A:200:CYS:HB3	1:A:207:ILE:HG23	1.99	0.45
1:B:244:ASN:HA	1:B:247:LYS:NZ	2.32	0.45
1:C:57:PRO:HG2	1:C:58:GLN:OE1	2.17	0.45
1:B:129:PRO:O	1:B:130:THR:HG23	2.17	0.44
1:C:80:VAL:HG21	1:D:25:TYR:CD1	2.52	0.44
1:D:8:VAL:HG21	1:D:224:ALA:HA	2.00	0.44
1:C:79:HIS:CE1	1:C:209:GLY:HA3	2.52	0.44
1:C:271:ALA:HB3	1:C:274:SER:OG	2.17	0.44
1:C:246:LYS:HD2	1:C:246:LYS:HA	1.62	0.44
1:D:303:ASP:HA	3:D:510:HOH:O	2.18	0.44
1:B:254:GLU:N	1:B:256:ILE:HD11	2.32	0.44
1:A:32:VAL:O	1:A:36:ALA:N	2.50	0.44
1:B:79:HIS:HB3	1:B:210:VAL:O	2.18	0.44
1:D:131:LYS:HA	1:D:132:PRO:HD3	1.65	0.44
1:D:295:ILE:C	1:D:296:THR:CG2	2.86	0.44
1:C:202:ARG:HA	1:C:206:VAL:O	2.18	0.44
1:C:4:GLY:HA3	1:C:46:GLY:O	2.18	0.44
1:D:199:PRO:HG2	1:D:212:ALA:O	2.17	0.44
1:C:18:GLN:HA	1:C:18:GLN:OE1	2.18	0.44
1:B:208:SER:HB3	3:B:555:HOH:O	2.18	0.43
1:C:129:PRO:HB2	1:C:165:VAL:O	2.18	0.43
1:A:308:LYS:HA	1:A:308:LYS:HD3	1.83	0.43
1:A:316:ASP:HA	3:A:1131:HOH:O	2.19	0.43
1:D:262:LEU:HD23	1:D:262:LEU:HA	1.74	0.43
1:A:248:PRO:C	1:A:250:GLU:H	2.21	0.43
1:D:10:ALA:HB2	1:D:220:THR:HG21	2.01	0.43
1:C:78:LEU:O	1:C:81:CYS:HB2	2.19	0.43
1:C:136:PHE:CE2	1:C:163:ARG:HD3	2.54	0.43
1:C:58:GLN:H	1:C:58:GLN:CD	2.21	0.43
1:B:242:CYS:SG	1:B:252:LYS:HG3	2.59	0.43
1:C:52:THR:CG2	1:C:53:SER:H	2.31	0.43
1:D:124:GLN:H	1:D:124:GLN:HG3	1.63	0.43
1:D:246:LYS:HD2	1:D:246:LYS:HA	1.45	0.43
1:A:78:LEU:HB3	1:A:211:ASP:OD2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:MET:O	1:B:97:MET:HB2	2.19	0.43
1:D:243:ILE:HG21	1:D:264:LEU:HD13	2.01	0.43
1:D:69:ALA:C	1:D:71:GLY:N	2.71	0.43
1:A:171:PRO:HG3	1:A:214:ILE:CG2	2.49	0.42
1:C:101:PHE:CZ	1:C:189:LEU:HD23	2.54	0.42
1:D:243:ILE:HB	1:D:269:HIS:ND1	2.33	0.42
1:C:63:LYS:HG2	1:C:133:VAL:HG13	2.01	0.42
1:D:166:VAL:HA	1:D:167:PRO:HD3	1.86	0.42
1:C:245:TYR:O	1:C:246:LYS:HB2	2.18	0.42
1:D:189:LEU:C	1:D:189:LEU:HD13	2.40	0.42
1:A:53:SER:C	1:A:216:LYS:HG3	2.39	0.42
1:A:87:GLY:HA3	1:B:83:ALA:O	2.20	0.42
1:C:120:ASP:OD1	1:C:122:LYS:N	2.52	0.42
1:D:170:ARG:HG3	1:D:171:PRO:HD2	2.02	0.42
1:D:18:GLN:NE2	3:D:534:HOH:O	2.52	0.42
1:D:233:MET:HE1	1:D:301:ALA:O	2.20	0.42
1:B:247:LYS:HG2	1:B:248:PRO:HD2	2.00	0.42
1:B:17:LEU:HD13	1:B:28:GLN:HG2	2.01	0.42
1:A:62:ILE:HD13	1:A:84:MET:HB2	2.00	0.42
1:B:274:SER:OG	2:B:401:ANP:H8	2.19	0.42
1:B:64:LEU:HD22	1:B:64:LEU:O	2.20	0.42
1:D:18:GLN:HB2	1:D:21:GLU:OE2	2.18	0.42
1:B:170:ARG:NH1	1:B:284:GLU:OE2	2.52	0.42
1:B:312:ARG:HD2	3:B:554:HOH:O	2.20	0.42
1:C:239:LEU:O	1:C:239:LEU:HD12	2.19	0.42
1:C:282:ALA:HB1	1:C:293:SER:CB	2.50	0.42
1:D:292:MET:HE3	1:D:314:ILE:HD11	2.00	0.42
1:A:33:GLU:HG2	3:A:1169:HOH:O	2.18	0.42
1:D:13:GLY:H	2:D:401:ANP:HNB1	1.66	0.42
1:D:78:LEU:HD23	1:D:212:ALA:N	2.35	0.42
1:C:305:LEU:HA	1:C:305:LEU:HD12	1.92	0.42
1:A:222:LEU:HA	1:A:225:LYS:HE3	2.01	0.42
1:B:263:ALA:HA	1:B:266:LYS:CG	2.50	0.42
1:D:168:SER:HA	1:D:213:VAL:O	2.20	0.42
1:D:189:LEU:HD22	1:D:189:LEU:HA	1.85	0.42
1:B:166:VAL:HB	1:B:167:PRO:CD	2.50	0.41
1:C:278:LYS:CE	2:C:401:ANP:O2B	2.68	0.41
1:A:11:LEU:O	1:A:216:LYS:HE3	2.21	0.41
1:D:255:GLU:C	1:D:256:ILE:HG13	2.40	0.41
1:A:84:MET:HG2	1:B:88:PHE:HB2	2.01	0.41
1:B:264:LEU:HD23	1:B:264:LEU:HA	1.80	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:247:LYS:O	1:D:249:ASP:N	2.52	0.41
1:A:84:MET:HB3	1:B:84:MET:HB3	2.02	0.41
1:C:69:ALA:C	1:C:71:GLY:H	2.24	0.41
1:D:11:LEU:HA	1:D:11:LEU:HD12	1.91	0.41
1:B:52:THR:HG21	1:B:220:THR:HG1	1.84	0.41
1:D:243:ILE:CG2	1:D:264:LEU:HD13	2.51	0.41
1:A:244:ASN:HB2	1:A:250:GLU:HA	2.02	0.41
1:B:62:ILE:HG22	1:B:81:CYS:SG	2.61	0.41
1:D:69:ALA:HB1	1:D:72:VAL:HG23	2.02	0.41
1:C:198:ILE:HA	1:C:199:PRO:HD3	1.96	0.41
1:A:97:MET:O	1:A:101:PHE:HD1	2.03	0.41
1:A:18:GLN:HB2	1:A:21:GLU:CD	2.42	0.41
1:D:186:ASN:ND2	3:D:529:HOH:O	2.31	0.41
1:A:198:ILE:HA	1:A:199:PRO:HD3	1.94	0.40
1:B:243:ILE:HG22	1:B:269:HIS:CB	2.51	0.40
1:A:110:CYS:HB2	1:A:189:LEU:O	2.22	0.40
1:A:258:LEU:HD22	1:A:286:THR:HG22	2.03	0.40
1:A:76:MET:HG3	1:A:84:MET:CE	2.52	0.40
1:D:1:MET:HE2	1:D:1:MET:HB3	1.98	0.40
1:D:227:LEU:HD12	1:D:227:LEU:HA	1.86	0.40
1:A:91:TYR:O	1:A:94:SER:HB2	2.21	0.40
1:B:110:CYS:HB3	3:B:541:HOH:O	2.21	0.40
1:B:170:ARG:NH2	1:B:218:LEU:HD13	2.36	0.40
1:C:256:ILE:HG12	1:C:257:LYS:N	2.36	0.40
1:D:103:ALA:HB2	3:D:583:HOH:O	2.22	0.40
1:C:143:LYS:O	1:C:143:LYS:HG2	2.21	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	281/317 (89%)	253 (90%)	24 (8%)	4 (1%)	11	22
1	B	281/317 (89%)	265 (94%)	13 (5%)	3 (1%)	14	30
1	C	314/317 (99%)	292 (93%)	19 (6%)	3 (1%)	15	32
1	D	314/317 (99%)	283 (90%)	27 (9%)	4 (1%)	12	24
All	All	1190/1268 (94%)	1093 (92%)	83 (7%)	14 (1%)	13	27

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	PRO
1	B	70	ALA
1	D	70	ALA
1	A	309	CYS
1	B	23	GLY
1	D	160	ARG
1	A	248	PRO
1	C	70	ALA
1	A	249	ASP
1	B	130	THR
1	C	309	CYS
1	D	246	LYS
1	D	248	PRO
1	C	23	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	234/257 (91%)	210 (90%)	24 (10%)	7	13
1	B	234/257 (91%)	199 (85%)	35 (15%)	3	5
1	C	257/257 (100%)	226 (88%)	31 (12%)	5	9
1	D	257/257 (100%)	215 (84%)	42 (16%)	2	3
All	All	982/1028 (96%)	850 (87%)	132 (13%)	4	6

All (132) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	11	LEU
1	A	17	LEU
1	A	20	LYS
1	A	37	SER
1	A	48	LYS
1	A	64	LEU
1	A	122	LYS
1	A	124	GLN
1	A	128	ASN
1	A	131	LYS
1	A	166	VAL
1	A	184	ILE
1	A	187	ASN
1	A	216	LYS
1	A	221	SER
1	A	227	LEU
1	A	239	LEU
1	A	247	LYS
1	A	250	GLU
1	A	257	LYS
1	A	266	LYS
1	A	284	GLU
1	A	296	THR
1	B	2	SER
1	B	7	VAL
1	B	11	LEU
1	B	17	LEU
1	B	20	LYS
1	B	22	LYS
1	B	29	ARG
1	B	59	VAL
1	B	63	LYS
1	B	64	LEU
1	B	72	VAL
1	B	73	SER
1	B	110	CYS
1	B	111	VAL
1	B	122	LYS
1	B	130	THR
1	B	165	VAL
1	B	168	SER

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Mol	Chain	Res	Type
1	B	184	ILE
1	B	189	LEU
1	B	193	THR
1	B	201	LYS
1	B	203	GLU
1	B	208	SER
1	B	216	LYS
1	B	227	LEU
1	B	239	LEU
1	B	247	LYS
1	B	252	LYS
1	B	253	LEU
1	B	256	ILE
1	B	266	LYS
1	B	292	MET
1	B	296	THR
1	B	305	LEU
1	C	2	SER
1	C	7	VAL
1	C	17	LEU
1	C	24	ASP
1	C	53	SER
1	C	63	LYS
1	C	64	LEU
1	C	66	ASN
1	C	94	SER
1	C	128	ASN
1	C	136	PHE
1	C	138	THR
1	C	141	GLU
1	C	149	ASN
1	C	153	ILE
1	C	189	LEU
1	C	193	THR
1	C	203	GLU
1	C	216	LYS
1	C	227	LEU
1	C	233	MET
1	C	238	VAL
1	C	246	LYS
1	C	252	LYS
1	C	253	LEU

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Mol	Chain	Res	Type
1	C	257	LYS
1	C	280	ARG
1	C	296	THR
1	C	302	VAL
1	C	305	LEU
1	C	308	LYS
1	D	2	SER
1	D	7	VAL
1	D	11	LEU
1	D	17	LEU
1	D	33	GLU
1	D	37	SER
1	D	53	SER
1	D	63	LYS
1	D	64	LEU
1	D	72	VAL
1	D	73	SER
1	D	94	SER
1	D	106	GLU
1	D	111	VAL
1	D	122	LYS
1	D	124	GLN
1	D	127	THR
1	D	135	ARG
1	D	141	GLU
1	D	144	ASP
1	D	149	ASN
1	D	152	LYS
1	D	154	LEU
1	D	156	GLU
1	D	163	ARG
1	D	165	VAL
1	D	170	ARG
1	D	172	LEU
1	D	193	THR
1	D	201	LYS
1	D	203	GLU
1	D	216	LYS
1	D	239	LEU
1	D	244	ASN
1	D	246	LYS
1	D	247	LYS

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Mol	Chain	Res	Type
1	D	249	ASP
1	D	256	ILE
1	D	265	GLU
1	D	266	LYS
1	D	296	THR
1	D	305	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	269	HIS

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 ⓘ

4 ligands are modelled in this entry.

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	$\# Z > 2$
2	ANP	D	401	-	29,33,33	2.13	8 (27%)	31,52,52	1.58	5 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ANP	A	1001	-	29,33,33	2.11	8 (27%)	31,52,52	1.74	7 (22%)
2	ANP	C	401	-	29,33,33	2.31	9 (31%)	31,52,52	1.72	8 (25%)
2	ANP	B	401	-	29,33,33	2.09	7 (24%)	31,52,52	1.68	7 (22%)

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
2	ANP	D	401	-	-	5/14/38/38	0/3/3/3
2	ANP	A	1001	-	-	5/14/38/38	0/3/3/3
2	ANP	C	401	-	-	4/14/38/38	0/3/3/3
2	ANP	B	401	-	-	4/14/38/38	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	ANP	PB-O1B	6.29	1.56	1.46
2	D	401	ANP	PB-O1B	5.87	1.55	1.46
2	A	1001	ANP	PB-O1B	5.83	1.55	1.46
2	B	401	ANP	PB-O1B	5.38	1.54	1.46
2	A	1001	ANP	C2'-C1'	-4.55	1.46	1.53
2	C	401	ANP	C2'-C1'	-4.47	1.47	1.53
2	B	401	ANP	C2'-C1'	-4.25	1.47	1.53
2	D	401	ANP	C2'-C1'	-3.97	1.47	1.53
2	C	401	ANP	PG-O1G	3.84	1.52	1.46
2	B	401	ANP	C6-N6	3.80	1.47	1.34
2	C	401	ANP	C6-N6	3.75	1.47	1.34
2	D	401	ANP	C6-N6	3.72	1.47	1.34
2	A	1001	ANP	C6-N6	3.60	1.47	1.34
2	D	401	ANP	O4'-C4'	-3.59	1.37	1.45
2	D	401	ANP	PG-O1G	3.52	1.51	1.46
2	C	401	ANP	O4'-C1'	3.46	1.45	1.41
2	B	401	ANP	PG-O1G	3.42	1.51	1.46
2	A	1001	ANP	C2'-C3'	-3.30	1.44	1.53
2	B	401	ANP	O4'-C4'	-3.27	1.37	1.45
2	A	1001	ANP	PG-O1G	3.26	1.51	1.46
2	A	1001	ANP	O4'-C4'	-3.24	1.37	1.45
2	C	401	ANP	C2'-C3'	-3.22	1.44	1.53
2	B	401	ANP	O4'-C1'	3.19	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ANP	C2'-C3'	-3.13	1.44	1.53
2	D	401	ANP	C2'-C3'	-3.06	1.45	1.53
2	C	401	ANP	PB-O3A	-3.04	1.55	1.59
2	C	401	ANP	O4'-C4'	-2.95	1.38	1.45
2	D	401	ANP	O4'-C1'	2.82	1.45	1.41
2	A	1001	ANP	O4'-C1'	2.80	1.45	1.41
2	C	401	ANP	PB-N3B	2.15	1.69	1.63
2	D	401	ANP	C2-N3	2.07	1.35	1.32
2	A	1001	ANP	O2'-C2'	-2.06	1.38	1.43

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	ANP	N3-C2-N1	-5.28	120.43	128.68
2	B	401	ANP	N3-C2-N1	-5.26	120.46	128.68
2	D	401	ANP	N3-C2-N1	-4.94	120.96	128.68
2	C	401	ANP	N3-C2-N1	-4.93	120.97	128.68
2	A	1001	ANP	C3'-C2'-C1'	3.81	106.71	100.98
2	B	401	ANP	C3'-C2'-C1'	3.66	106.48	100.98
2	D	401	ANP	C3'-C2'-C1'	3.37	106.06	100.98
2	C	401	ANP	PA-O3A-PB	-3.27	121.11	132.62
2	C	401	ANP	C3'-C2'-C1'	3.22	105.83	100.98
2	A	1001	ANP	O3A-PB-N3B	3.21	115.50	106.59
2	C	401	ANP	O5'-C5'-C4'	3.20	120.00	108.99
2	C	401	ANP	O3A-PB-N3B	3.01	114.94	106.59
2	A	1001	ANP	PA-O3A-PB	-2.86	122.55	132.62
2	B	401	ANP	O3A-PB-N3B	2.69	114.05	106.59
2	D	401	ANP	PA-O3A-PB	-2.40	124.17	132.62
2	B	401	ANP	C2-N1-C6	2.38	122.82	118.75
2	A	1001	ANP	C2-N1-C6	2.34	122.75	118.75
2	B	401	ANP	PA-O3A-PB	-2.32	124.45	132.62
2	D	401	ANP	C2-N1-C6	2.31	122.71	118.75
2	D	401	ANP	C1'-N9-C4	-2.31	122.59	126.64
2	C	401	ANP	C2-N1-C6	2.21	122.54	118.75
2	B	401	ANP	O5'-C5'-C4'	2.16	116.41	108.99
2	B	401	ANP	C1'-N9-C4	-2.15	122.87	126.64
2	A	1001	ANP	O5'-C5'-C4'	2.06	116.08	108.99
2	C	401	ANP	O2G-PG-O1G	-2.06	108.27	113.45
2	C	401	ANP	O3G-PG-O1G	-2.02	108.38	113.45
2	A	1001	ANP	O2G-PG-O1G	-2.01	108.40	113.45

There are no chirality outliers.

All (18) torsion outliers are listed below:

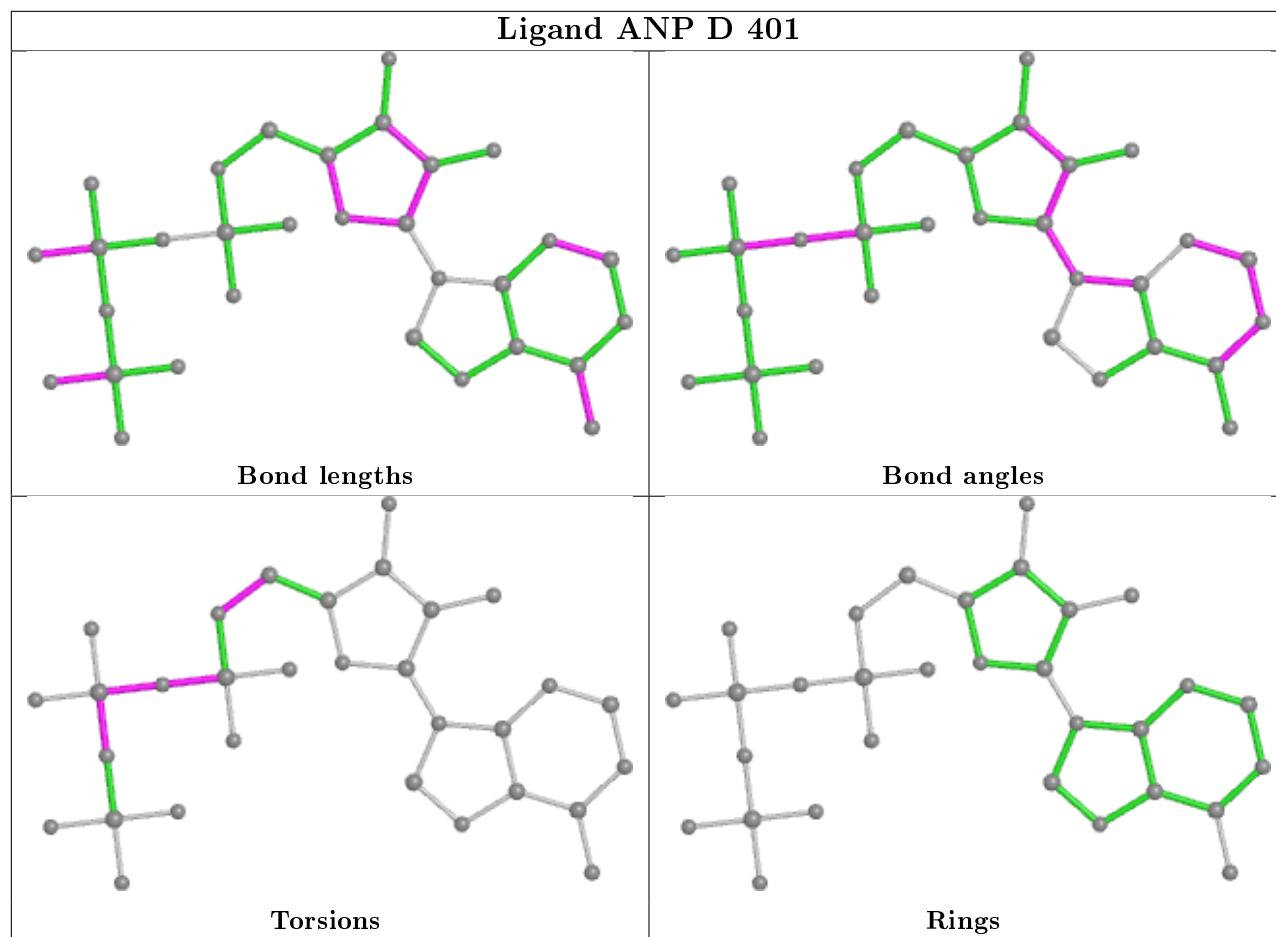
Mol	Chain	Res	Type	Atoms
2	D	401	ANP	PG-N3B-PB-O1B
2	D	401	ANP	PA-O3A-PB-O1B
2	D	401	ANP	PA-O3A-PB-O2B
2	A	1001	ANP	PG-N3B-PB-O1B
2	A	1001	ANP	PA-O3A-PB-O1B
2	A	1001	ANP	PA-O3A-PB-O2B
2	C	401	ANP	PG-N3B-PB-O1B
2	C	401	ANP	PA-O3A-PB-O1B
2	B	401	ANP	PG-N3B-PB-O1B
2	B	401	ANP	PA-O3A-PB-O1B
2	B	401	ANP	PA-O3A-PB-O2B
2	B	401	ANP	C4'-C5'-O5'-PA
2	D	401	ANP	PB-O3A-PA-O1A
2	A	1001	ANP	PB-O3A-PA-O1A
2	D	401	ANP	C4'-C5'-O5'-PA
2	C	401	ANP	C4'-C5'-O5'-PA
2	A	1001	ANP	C4'-C5'-O5'-PA
2	C	401	ANP	PB-O3A-PA-O1A

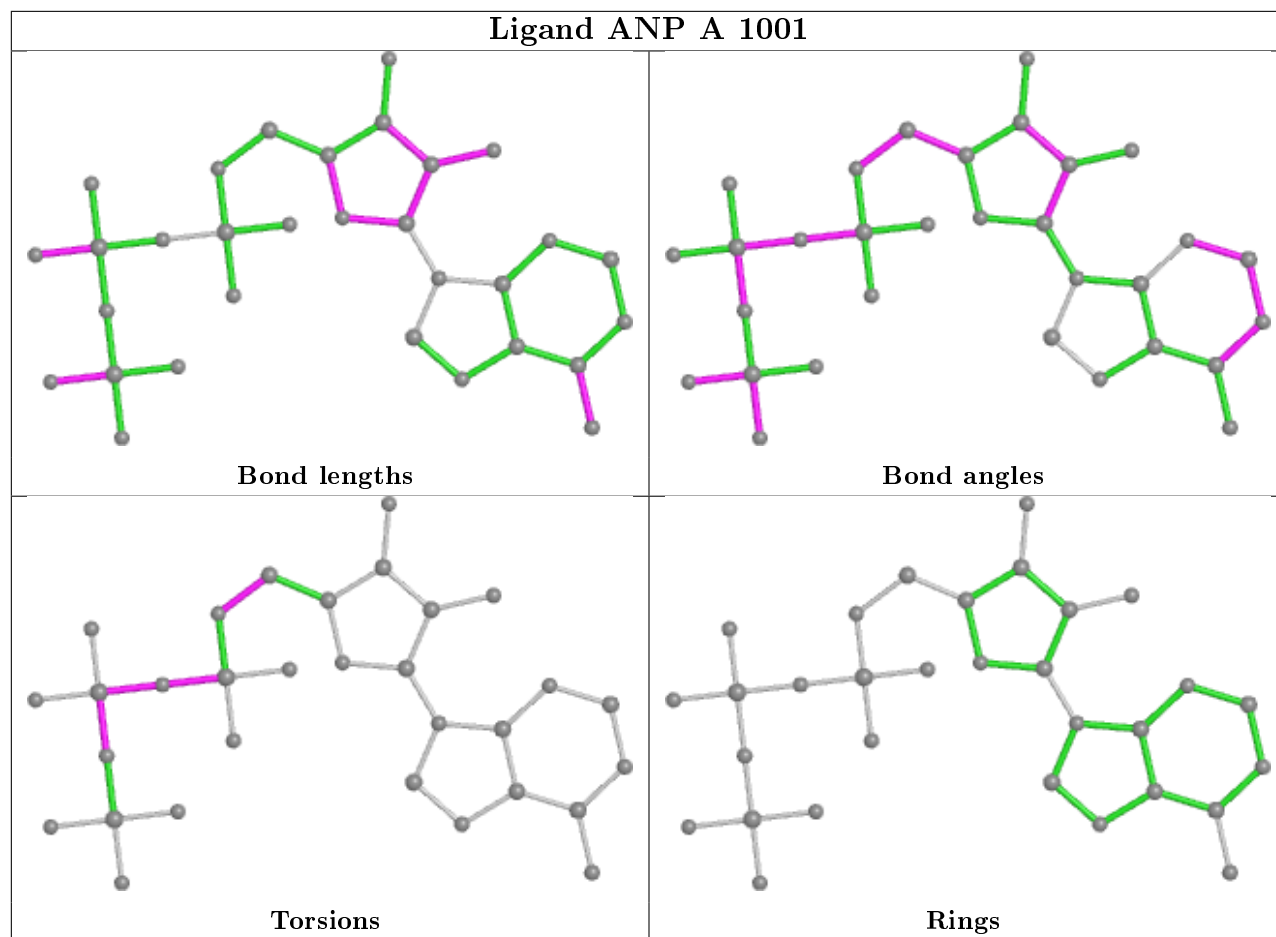
There are no ring outliers.

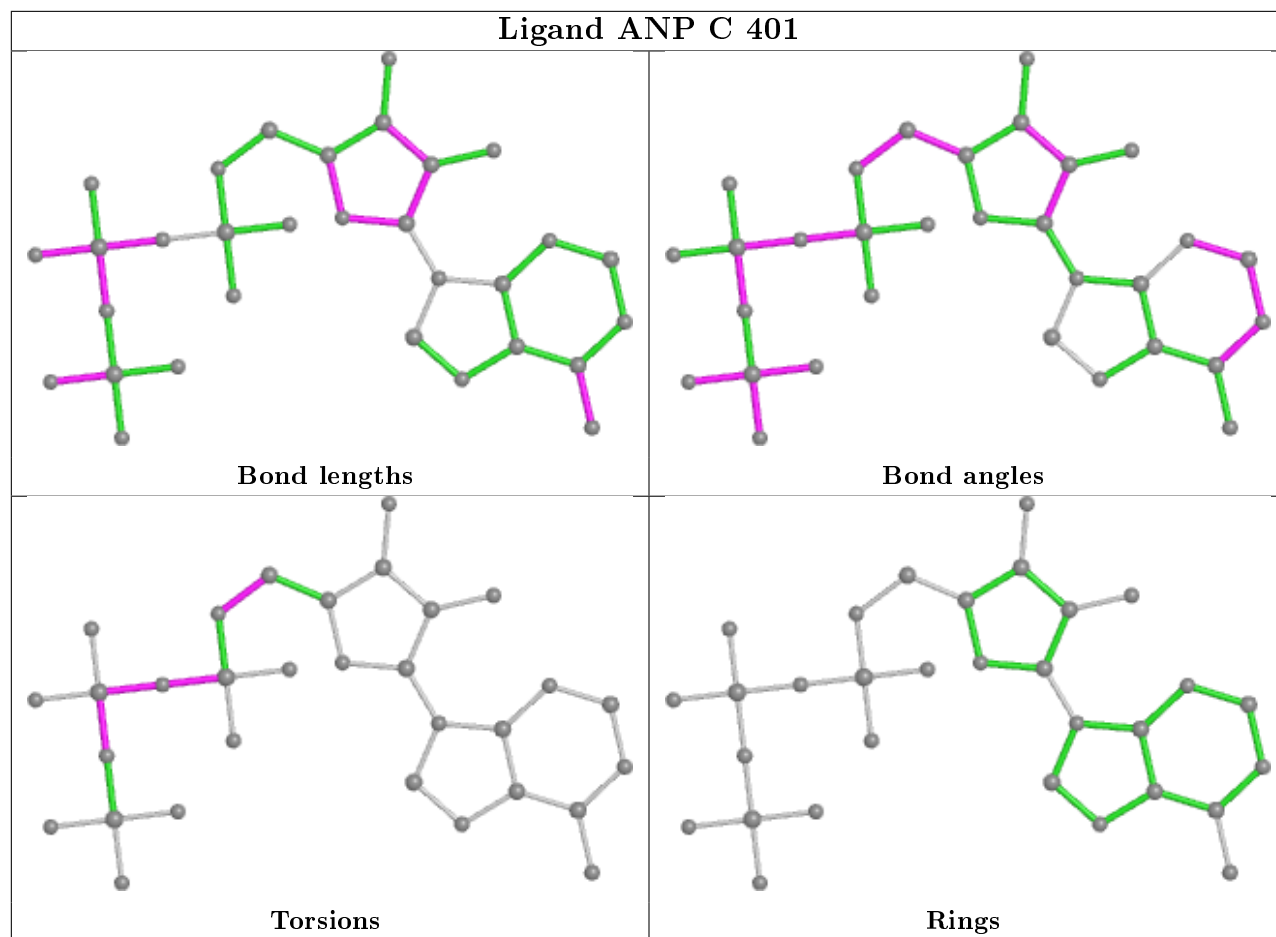
4 monomers are involved in 16 short contacts:

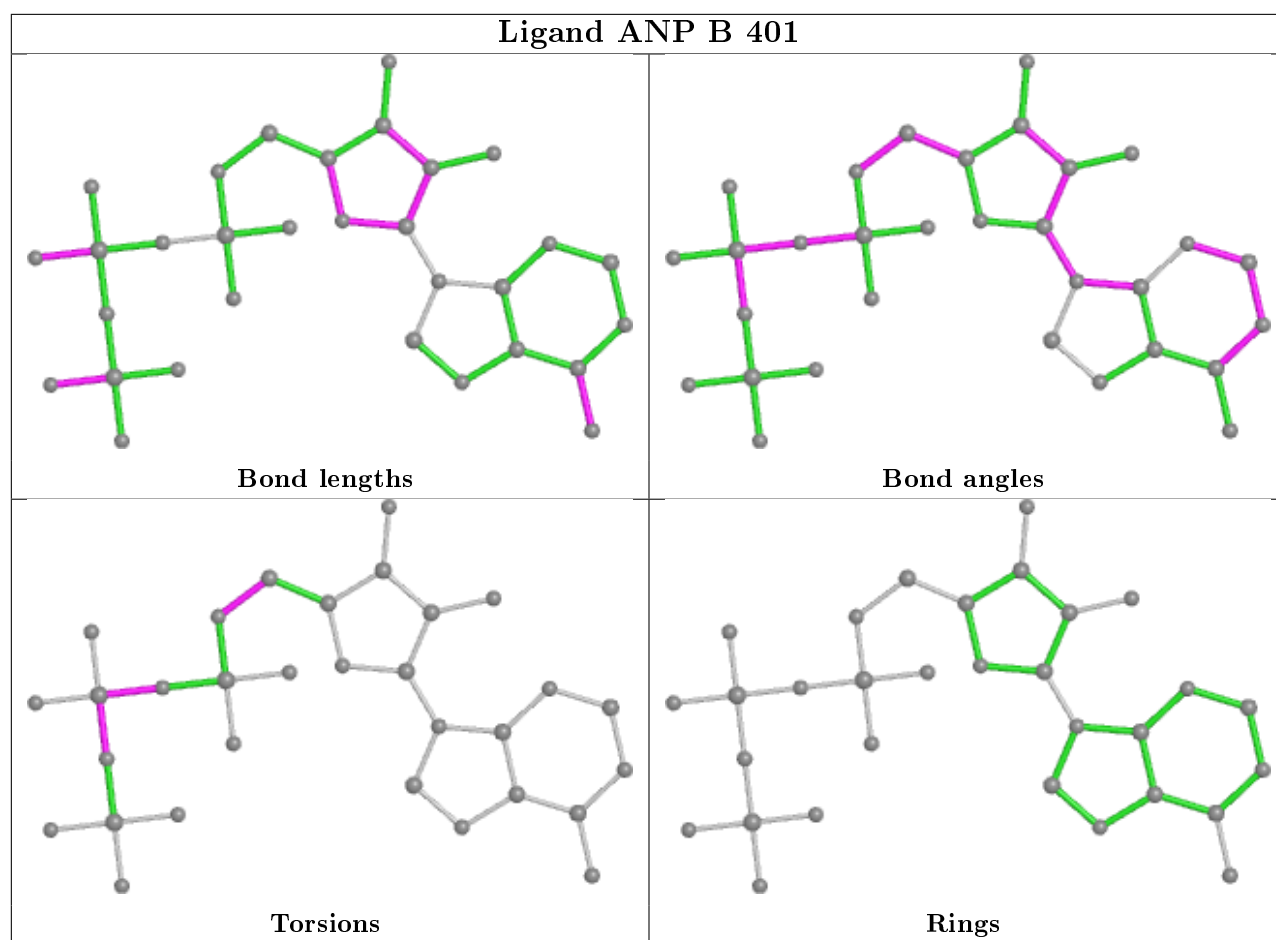
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	ANP	4	0
2	A	1001	ANP	1	0
2	C	401	ANP	5	0
2	B	401	ANP	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









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	285/317 (89%)	-0.59	1 (0%) 92 91	5, 16, 44, 68	0
1	B	285/317 (89%)	-0.51	2 (0%) 87 86	6, 18, 48, 68	0
1	C	316/317 (99%)	-0.31	10 (3%) 47 40	6, 20, 70, 94	0
1	D	316/317 (99%)	-0.33	13 (4%) 37 30	6, 20, 70, 90	0
All	All	1202/1268 (94%)	-0.43	26 (2%) 62 56	5, 18, 60, 94	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	148	ALA	8.5
1	B	165	VAL	5.7
1	D	142	ALA	5.3
1	D	151	GLY	4.9
1	D	150	PRO	4.8
1	C	144	ASP	4.6
1	A	165	VAL	4.4
1	C	150	PRO	3.6
1	D	144	ASP	3.5
1	D	140	GLN	3.5
1	C	146	MET	3.1
1	C	145	LEU	2.9
1	D	158	ALA	2.5
1	D	157	ASP	2.5
1	D	159	GLY	2.4
1	C	142	ALA	2.4
1	D	133	VAL	2.4
1	D	134	GLY	2.3
1	B	248	PRO	2.3
1	C	164	VAL	2.3
1	D	19	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	143	LYS	2.2
1	D	149	ASN	2.2
1	C	157	ASP	2.1
1	C	135	ARG	2.1
1	C	133	VAL	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](#)

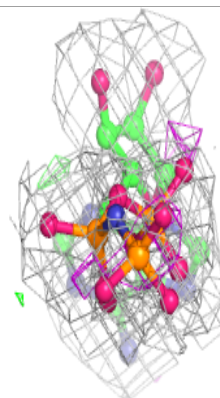
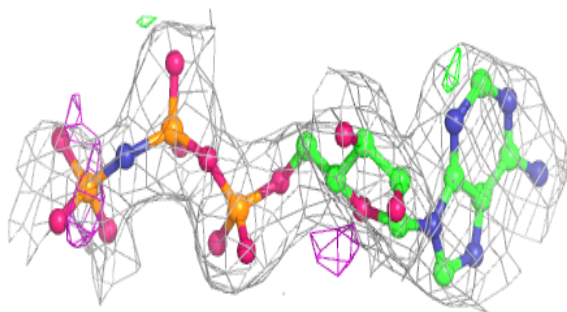
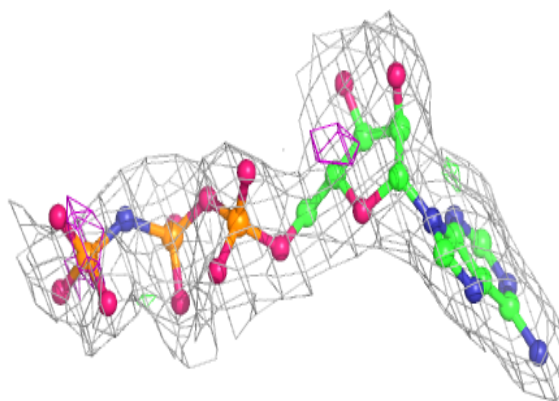
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ANP	D	401	31/31	0.93	0.16	17,29,49,81	0
2	ANP	C	401	31/31	0.93	0.15	15,29,42,77	0
2	ANP	B	401	31/31	0.93	0.17	15,28,38,82	0
2	ANP	A	1001	31/31	0.94	0.14	16,24,44,76	0

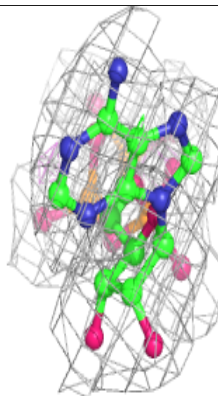
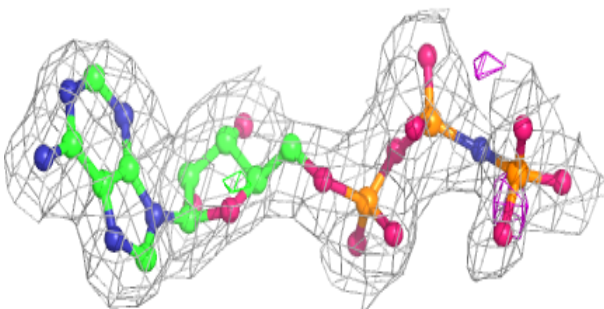
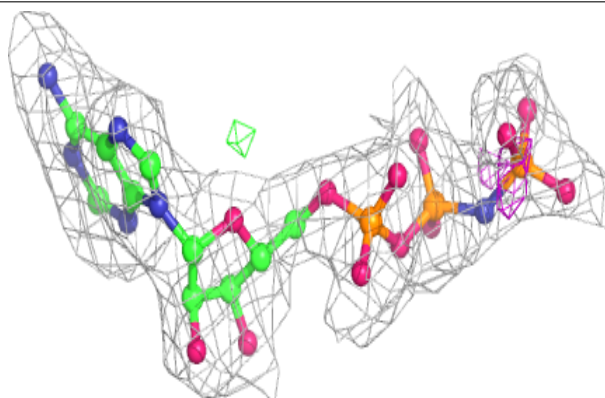
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ANP D 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

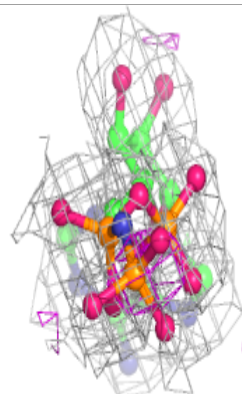
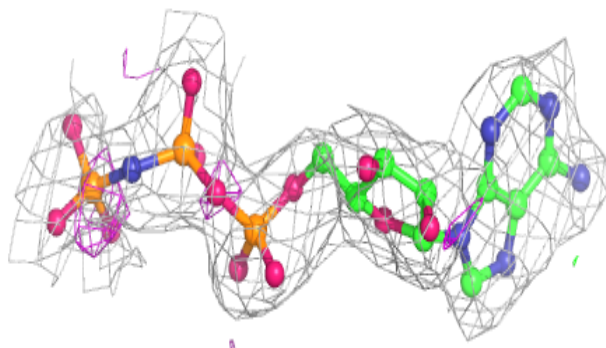
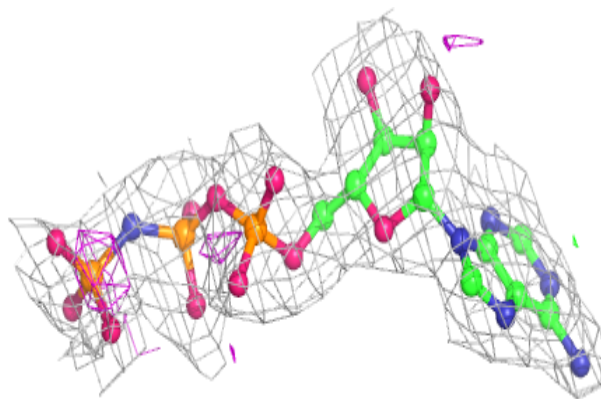
**Electron density around ANP C 401:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

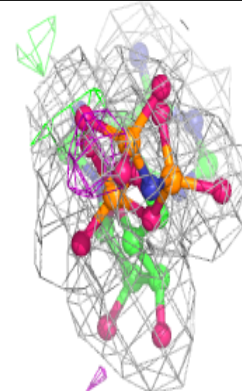
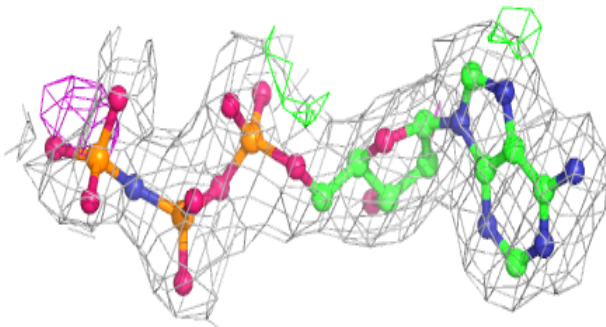
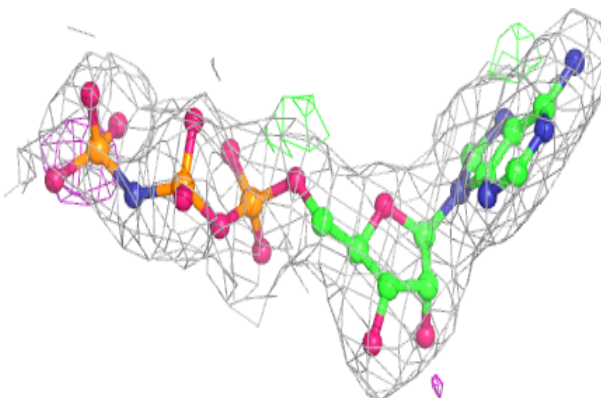


Electron density around ANP B 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around ANP A 1001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.