



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 08:51 am BST

PDB ID : 3H39
Title : The complex structure of CCA-adding enzyme with ATP
Authors : Toh, Y.; Tomita, K.
Deposited on : 2009-04-16
Resolution : 2.85 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

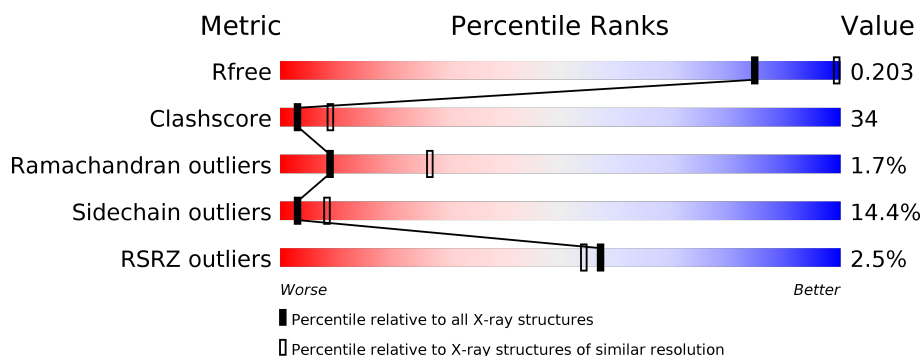
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	441	<div> <div>3%</div> <div> <div></div> <div>50%</div> <div>34%</div> <div>9%</div> <div>6%</div> </div> </div>
1	B	441	<div> <div>2%</div> <div> <div></div> <div>46%</div> <div>40%</div> <div>7%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6970 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 TRNA nucleotidyl transferase-related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3418	2208	580	619	11			
1	B	420	Total	C	N	O	S	0	0	0
			3450	2228	586	625	11			

There are 28 discrepancies between the modelled and reference sequences:

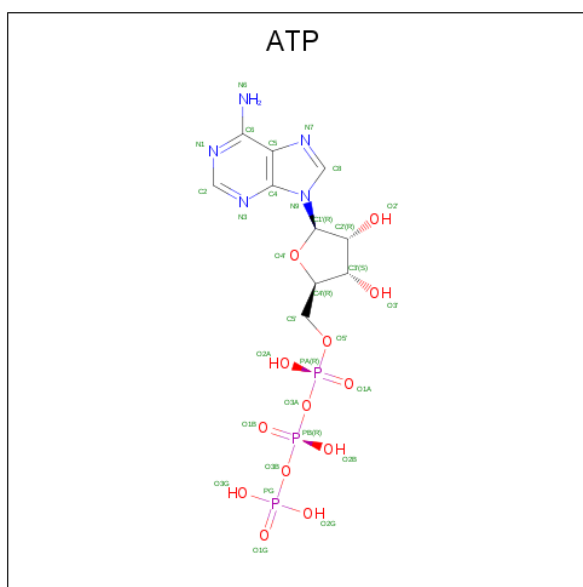
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q9WZH4
A	429	LYS	-	EXPRESSION TAG	UNP Q9WZH4
A	430	LEU	-	EXPRESSION TAG	UNP Q9WZH4
A	431	ALA	-	EXPRESSION TAG	UNP Q9WZH4
A	432	ALA	-	EXPRESSION TAG	UNP Q9WZH4
A	433	ALA	-	EXPRESSION TAG	UNP Q9WZH4
A	434	LEU	-	EXPRESSION TAG	UNP Q9WZH4
A	435	GLU	-	EXPRESSION TAG	UNP Q9WZH4
A	436	HIS	-	EXPRESSION TAG	UNP Q9WZH4
A	437	HIS	-	EXPRESSION TAG	UNP Q9WZH4
A	438	HIS	-	EXPRESSION TAG	UNP Q9WZH4
A	439	HIS	-	EXPRESSION TAG	UNP Q9WZH4
A	440	HIS	-	EXPRESSION TAG	UNP Q9WZH4
A	441	HIS	-	EXPRESSION TAG	UNP Q9WZH4
B	1	MET	-	EXPRESSION TAG	UNP Q9WZH4
B	429	LYS	-	EXPRESSION TAG	UNP Q9WZH4
B	430	LEU	-	EXPRESSION TAG	UNP Q9WZH4
B	431	ALA	-	EXPRESSION TAG	UNP Q9WZH4
B	432	ALA	-	EXPRESSION TAG	UNP Q9WZH4
B	433	ALA	-	EXPRESSION TAG	UNP Q9WZH4
B	434	LEU	-	EXPRESSION TAG	UNP Q9WZH4
B	435	GLU	-	EXPRESSION TAG	UNP Q9WZH4
B	436	HIS	-	EXPRESSION TAG	UNP Q9WZH4
B	437	HIS	-	EXPRESSION TAG	UNP Q9WZH4
B	438	HIS	-	EXPRESSION TAG	UNP Q9WZH4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	439	HIS	-	EXPRESSION TAG	UNP Q9WZH4
B	440	HIS	-	EXPRESSION TAG	UNP Q9WZH4
B	441	HIS	-	EXPRESSION TAG	UNP Q9WZH4

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

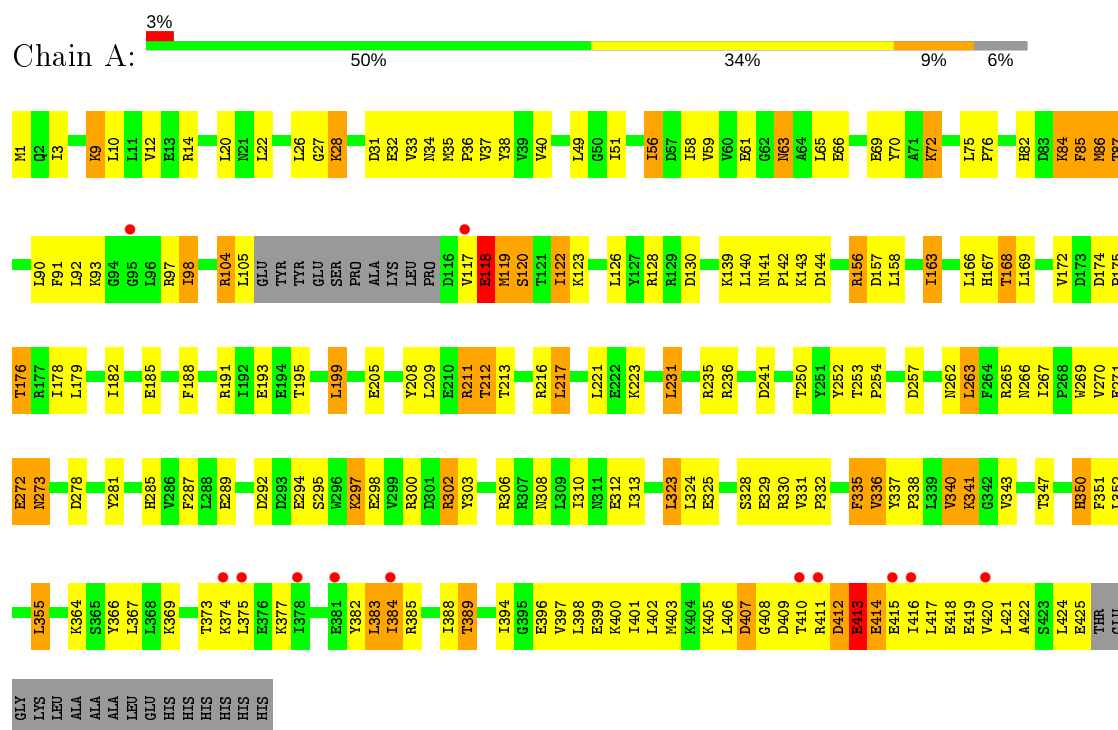
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		
3	B	10	Total	O	0	0
			10	10		

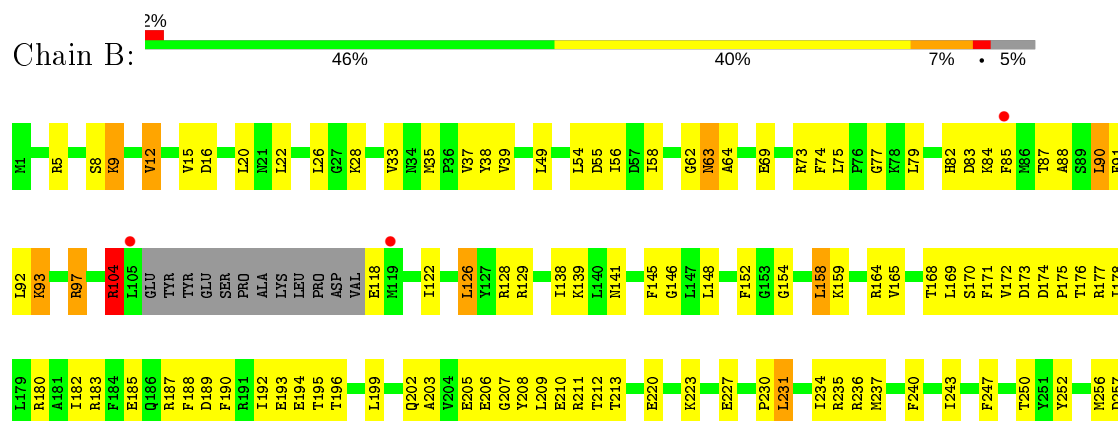
3 Residue-property plots

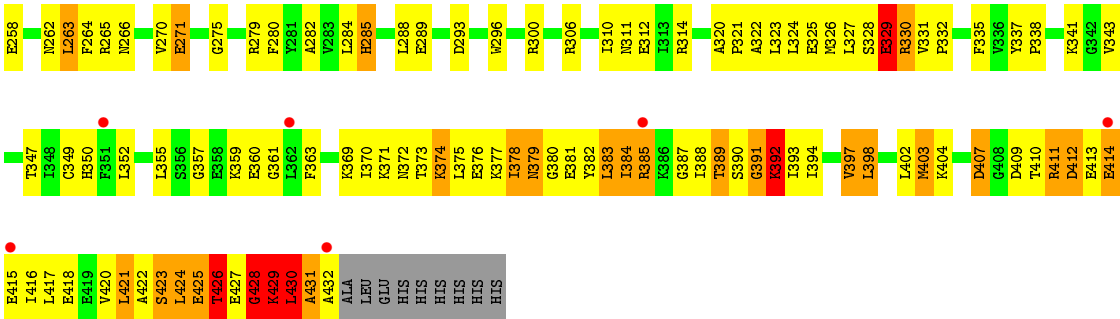
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: TRNA nucleotidyl transferase-related protein



- Molecule 1: TRNA nucleotidyl transferase-related protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	186.09Å 64.07Å 151.77Å 90.00° 100.88° 90.00°	Depositor
Resolution (Å)	36.03 – 2.85 48.46 – 2.85	Depositor EDS
% Data completeness (in resolution range)	96.9 (36.03-2.85) 96.8 (48.46-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.210 , 0.252 0.207 , 0.203	Depositor DCC
R_{free} test set	1963 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	63.7	Xtriage
Anisotropy	0.171	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 62.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6970	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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: ATP

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	0.66	0/3479	0.79	5/4674 (0.1%)
1	B	0.49	0/3511	0.77	7/4716 (0.1%)
All	All	0.58	0/6990	0.78	12/9390 (0.1%)

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	B	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	104	ARG	CB-CA-C	-17.04	76.31	110.40
1	B	392	LYS	CB-CA-C	-11.00	88.41	110.40
1	B	430	LEU	CB-CA-C	10.41	129.98	110.20
1	B	104	ARG	C-N-CA	6.68	138.40	121.70
1	B	425	GLU	N-CA-C	-6.19	94.29	111.00
1	B	428	GLY	N-CA-C	5.99	128.08	113.10
1	A	413	GLU	CB-CA-C	-5.71	98.99	110.40
1	A	188	PHE	N-CA-C	-5.67	95.70	111.00
1	A	120	SER	CB-CA-C	5.32	120.20	110.10
1	B	387	GLY	N-CA-C	-5.28	99.91	113.10
1	A	341	LYS	CB-CA-C	-5.23	99.94	110.40
1	A	272	GLU	CB-CA-C	-5.07	100.26	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	104	ARG	Peptide

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	3418	0	3517	232	0
1	B	3450	0	3554	243	0
2	A	31	0	12	1	0
2	B	31	0	12	3	0
3	A	30	0	0	0	0
3	B	10	0	0	1	0
All	All	6970	0	7095	476	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 34.

All (476) 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:90:LEU:HD11	1:A:92:LEU:CD2	1.62	1.27
1:B:427:GLU:CG	1:B:428:GLY:H	1.47	1.20
1:B:424:LEU:HD21	1:B:431:ALA:CB	1.71	1.19
1:B:104:ARG:HD3	1:B:128:ARG:HD3	1.19	1.18
1:B:381:GLU:HG3	1:B:382:TYR:H	1.01	1.18
1:A:383:LEU:CD1	1:A:383:LEU:H	1.52	1.16
1:B:202:GLN:O	1:B:206:GLU:HG3	1.42	1.16
1:A:383:LEU:HD12	1:A:383:LEU:N	1.57	1.13
1:A:410:THR:HG22	1:A:411:ARG:N	1.57	1.13
1:B:425:GLU:O	1:B:426:THR:HG22	1.48	1.13
1:A:341:LYS:O	1:A:341:LYS:HG3	1.50	1.11
1:A:410:THR:CG2	1:A:411:ARG:H	1.64	1.11
1:B:411:ARG:HD2	1:B:411:ARG:O	1.51	1.09
1:A:90:LEU:HD11	1:A:92:LEU:HD21	1.12	1.09
1:B:424:LEU:CD2	1:B:431:ALA:HB1	1.83	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:GLU:HG2	1:B:428:GLY:H	0.93	1.08
1:A:90:LEU:CD1	1:A:92:LEU:CD2	2.29	1.07
1:B:430:LEU:HD12	1:B:431:ALA:H	1.17	1.07
1:B:429:LYS:O	1:B:430:LEU:HB3	1.52	1.04
1:B:330:ARG:HG2	1:B:330:ARG:O	1.58	1.04
1:A:410:THR:HG22	1:A:411:ARG:H	0.92	1.03
1:B:427:GLU:CG	1:B:428:GLY:N	2.13	1.02
1:A:90:LEU:CD1	1:A:92:LEU:HD21	1.89	1.01
1:A:331:VAL:HB	1:A:332:PRO:HD2	1.39	1.00
1:A:82:HIS:HD2	1:A:87:THR:CG2	1.74	0.99
1:A:117:VAL:O	1:A:118:GLU:HB3	1.59	0.99
1:A:273:ASN:N	1:A:273:ASN:HD22	1.60	0.98
1:A:117:VAL:O	1:A:118:GLU:CB	2.10	0.98
1:B:427:GLU:HG2	1:B:428:GLY:N	1.69	0.98
1:A:384:ILE:HD11	1:A:397:VAL:HG21	1.45	0.97
1:B:424:LEU:HD21	1:B:431:ALA:HB1	0.97	0.97
1:A:383:LEU:H	1:A:383:LEU:HD12	0.82	0.96
1:A:373:THR:O	1:A:373:THR:HG22	1.66	0.95
1:B:104:ARG:HD3	1:B:128:ARG:CD	1.96	0.95
1:B:381:GLU:HG3	1:B:382:TYR:N	1.82	0.93
1:B:176:THR:HG23	1:B:220:GLU:HG3	1.51	0.92
1:B:388:ILE:O	1:B:389:THR:HG23	1.69	0.92
1:B:381:GLU:CG	1:B:382:TYR:H	1.76	0.91
1:B:388:ILE:O	1:B:389:THR:OG1	1.88	0.91
1:B:328:SER:O	1:B:329:GLU:HB2	1.66	0.91
1:A:82:HIS:CD2	1:A:87:THR:CG2	2.55	0.90
1:A:90:LEU:HD11	1:A:92:LEU:HD23	1.53	0.89
1:A:82:HIS:HD2	1:A:87:THR:HG23	1.37	0.88
1:B:411:ARG:C	1:B:411:ARG:HD2	1.89	0.86
1:B:381:GLU:O	1:B:384:ILE:HG23	1.77	0.85
1:B:429:LYS:O	1:B:430:LEU:CB	2.24	0.85
1:B:93:LYS:HD2	1:B:93:LYS:H	1.40	0.85
1:A:90:LEU:CD1	1:A:92:LEU:HD23	2.07	0.85
1:A:9:LYS:CD	1:A:9:LYS:H	1.90	0.85
1:A:37:VAL:H	1:A:141:ASN:HD21	1.24	0.84
1:B:37:VAL:H	1:B:141:ASN:HD21	1.26	0.84
1:A:272:GLU:HB2	1:A:273:ASN:ND2	1.93	0.84
1:B:430:LEU:HD12	1:B:431:ALA:N	1.93	0.83
1:B:425:GLU:C	1:B:426:THR:HG22	1.96	0.83
1:B:69:GLU:OE1	1:B:79:LEU:CD2	2.28	0.81
1:A:417:LEU:HA	1:A:420:VAL:HG22	1.61	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:ASP:OD1	1:B:176:THR:HB	1.80	0.80
1:B:388:ILE:O	1:B:389:THR:CG2	2.29	0.80
1:A:273:ASN:N	1:A:273:ASN:ND2	2.30	0.79
1:A:297:LYS:HG3	1:A:298:GLU:N	1.97	0.79
1:A:93:LYS:H	1:A:93:LYS:HD2	1.46	0.79
1:A:139:LYS:HD3	1:A:144:ASP:HB2	1.64	0.79
1:A:373:THR:CG2	1:A:373:THR:O	2.30	0.79
1:B:411:ARG:CD	1:B:411:ARG:O	2.30	0.79
1:A:91:PHE:HZ	1:A:97:ARG:HH11	1.27	0.79
1:B:420:VAL:O	1:B:423:SER:HB2	1.82	0.78
1:B:425:GLU:O	1:B:426:THR:CG2	2.29	0.78
1:A:49:LEU:HD21	1:A:158:LEU:HD22	1.64	0.78
1:B:330:ARG:CG	1:B:330:ARG:O	2.32	0.78
1:B:427:GLU:HG3	1:B:428:GLY:N	1.98	0.77
1:A:118:GLU:C	1:A:120:SER:H	1.87	0.77
1:B:93:LYS:CD	1:B:93:LYS:H	1.97	0.77
1:A:262:ASN:HD22	1:A:265:ARG:HH22	1.31	0.77
1:B:331:VAL:HG12	1:B:332:PRO:HD2	1.67	0.76
1:B:380:GLY:O	1:B:381:GLU:HG2	1.84	0.76
1:B:12:VAL:HA	1:B:20:LEU:HD11	1.66	0.76
1:A:56:ILE:O	1:A:56:ILE:HG12	1.85	0.76
1:B:328:SER:O	1:B:329:GLU:CB	2.35	0.75
1:B:424:LEU:N	1:B:424:LEU:HD13	2.02	0.75
1:A:384:ILE:HG21	1:A:394:ILE:HG23	1.68	0.74
1:A:90:LEU:HB3	1:A:98:ILE:CG2	2.17	0.74
1:A:294:GLU:OE1	1:A:297:LYS:HE2	1.87	0.74
1:B:410:THR:HG21	1:B:414:GLU:HG2	1.70	0.74
1:B:209:LEU:H	1:B:209:LEU:HD12	1.52	0.74
1:B:388:ILE:HG22	1:B:389:THR:N	2.03	0.74
1:B:413:GLU:O	1:B:416:ILE:HG22	1.88	0.74
1:A:117:VAL:O	1:A:118:GLU:CG	2.36	0.73
1:A:341:LYS:CG	1:A:341:LYS:O	2.30	0.73
1:A:211:ARG:HH21	1:A:211:ARG:CG	2.01	0.73
1:B:411:ARG:C	1:B:411:ARG:CD	2.52	0.72
1:B:388:ILE:O	1:B:389:THR:CB	2.37	0.72
1:B:380:GLY:O	1:B:381:GLU:CG	2.37	0.72
1:B:247:PHE:O	1:B:250:THR:HG22	1.89	0.72
1:A:328:SER:O	1:A:329:GLU:CB	2.34	0.72
1:A:82:HIS:CD2	1:A:87:THR:HG22	2.24	0.72
1:A:117:VAL:O	1:A:118:GLU:HG2	1.89	0.72
1:A:82:HIS:HB2	1:A:87:THR:HG22	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:427:GLU:HG3	1:B:428:GLY:H	1.48	0.72
1:A:410:THR:HG21	1:A:414:GLU:HB3	1.72	0.72
1:B:424:LEU:N	1:B:424:LEU:CD1	2.53	0.72
1:A:340:VAL:HG23	1:A:340:VAL:O	1.89	0.71
1:A:355:LEU:HD22	1:A:355:LEU:H	1.56	0.71
1:B:388:ILE:CG2	1:B:389:THR:H	2.02	0.71
1:A:424:LEU:O	1:A:425:GLU:HG2	1.91	0.70
1:B:428:GLY:O	1:B:429:LYS:HB3	1.92	0.70
1:B:194:GLU:HG3	1:B:195:THR:N	2.05	0.70
1:B:202:GLN:HG2	1:B:206:GLU:OE2	1.92	0.69
1:B:388:ILE:CG2	1:B:389:THR:N	2.55	0.69
1:A:9:LYS:HD2	1:A:9:LYS:H	1.57	0.69
1:B:262:ASN:HD22	1:B:265:ARG:HH12	1.41	0.69
1:B:74:PHE:C	1:B:75:LEU:HD12	2.13	0.69
1:B:54:LEU:O	1:B:56:ILE:HD12	1.92	0.69
1:B:56:ILE:CG2	1:B:58:ILE:HD12	2.23	0.69
1:B:56:ILE:HG22	1:B:56:ILE:O	1.91	0.69
1:B:409:ASP:OD1	1:B:409:ASP:C	2.29	0.68
1:B:376:GLU:OE1	1:B:376:GLU:HA	1.91	0.68
1:A:174:ASP:OD1	1:A:176:THR:HB	1.94	0.68
1:A:410:THR:CG2	1:A:414:GLU:HB3	2.23	0.68
1:A:420:VAL:O	1:A:424:LEU:HB2	1.94	0.68
1:A:388:ILE:O	1:A:388:ILE:HG23	1.94	0.67
1:A:85:PHE:O	1:A:87:THR:HB	1.95	0.67
1:B:410:THR:HG22	1:B:411:ARG:H	1.58	0.67
1:B:175:PRO:HD2	1:B:212:THR:HG21	1.75	0.67
1:B:384:ILE:HD11	1:B:394:ILE:HA	1.77	0.67
1:B:407:ASP:OD2	1:B:407:ASP:N	2.26	0.67
1:A:384:ILE:HG23	1:A:394:ILE:HD13	1.77	0.66
1:B:428:GLY:O	1:B:429:LYS:CB	2.44	0.66
1:A:84:LYS:O	1:A:84:LYS:HD2	1.95	0.66
1:B:380:GLY:C	1:B:381:GLU:HG2	2.15	0.66
1:B:378:ILE:HG21	1:B:411:ARG:HH21	1.59	0.66
1:A:213:THR:OG1	1:A:216:ARG:HG3	1.95	0.65
1:A:84:LYS:C	1:A:84:LYS:HD2	2.17	0.65
1:A:328:SER:O	1:A:329:GLU:HB3	1.97	0.64
1:A:85:PHE:O	1:A:86:MET:C	2.35	0.64
1:A:117:VAL:HG23	1:A:118:GLU:HG3	1.80	0.64
1:A:252:TYR:CE2	1:A:257:ASP:HB2	2.32	0.64
1:A:403:MET:HA	1:A:406:LEU:HD12	1.79	0.63
1:B:381:GLU:CG	1:B:382:TYR:N	2.49	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LEU:HD13	1:A:91:PHE:N	2.13	0.63
1:A:411:ARG:O	1:A:413:GLU:N	2.32	0.63
1:A:90:LEU:HD13	1:A:90:LEU:C	2.17	0.63
1:B:172:VAL:HG13	1:B:208:TYR:CE1	2.34	0.62
1:A:117:VAL:C	1:A:118:GLU:CG	2.65	0.62
1:B:62:GLY:O	1:B:63:ASN:HB2	2.00	0.62
1:B:252:TYR:HA	1:B:256:MET:HG2	1.82	0.62
1:A:262:ASN:ND2	1:A:265:ARG:HH12	1.97	0.62
1:A:272:GLU:HB2	1:A:273:ASN:HD22	1.62	0.62
1:B:391:GLY:O	1:B:392:LYS:HB2	1.99	0.61
1:A:331:VAL:HB	1:A:332:PRO:CD	2.17	0.61
1:B:331:VAL:CG1	1:B:332:PRO:HD2	2.30	0.61
1:A:1:MET:SD	1:A:156:ARG:HD2	2.41	0.60
1:A:323:LEU:HD12	1:A:351:PHE:CD2	2.35	0.60
1:A:302:ARG:HG2	1:A:303:TYR:CE1	2.35	0.60
1:B:410:THR:HG22	1:B:411:ARG:N	2.15	0.60
1:B:69:GLU:OE1	1:B:79:LEU:HD23	2.01	0.60
1:A:384:ILE:CG2	1:A:394:ILE:HD13	2.32	0.60
1:B:194:GLU:CG	1:B:195:THR:N	2.64	0.60
1:B:329:GLU:O	1:B:331:VAL:N	2.34	0.60
1:B:337:TYR:N	1:B:338:PRO:HD2	2.17	0.60
1:A:211:ARG:HG2	1:A:211:ARG:HH21	1.66	0.60
1:A:407:ASP:OD2	1:A:407:ASP:N	2.30	0.60
1:A:297:LYS:HG3	1:A:298:GLU:H	1.67	0.59
1:B:296:TRP:CE2	1:B:310:ILE:HB	2.37	0.59
1:A:63:ASN:HD21	1:A:65:LEU:HB2	1.67	0.59
1:A:355:LEU:HD22	1:A:355:LEU:N	2.15	0.59
1:A:93:LYS:N	1:A:93:LYS:HD2	2.14	0.59
1:B:285:HIS:HB3	1:B:350:HIS:CD2	2.37	0.59
1:A:417:LEU:HA	1:A:420:VAL:CG2	2.31	0.59
1:B:49:LEU:HD21	1:B:158:LEU:HD22	1.84	0.59
1:B:388:ILE:C	1:B:389:THR:OG1	2.39	0.59
1:A:383:LEU:CD1	1:A:383:LEU:N	2.30	0.59
1:B:393:ILE:O	1:B:397:VAL:HG13	2.03	0.59
1:A:384:ILE:O	1:A:389:THR:OG1	2.21	0.58
1:A:130:ASP:HB3	2:A:501:ATP:C2	2.38	0.58
1:B:262:ASN:ND2	1:B:265:ARG:HH12	2.01	0.58
1:B:263:LEU:HD11	1:B:282:ALA:O	2.02	0.58
1:A:120:SER:HA	1:A:122:ILE:CD1	2.34	0.58
1:A:118:GLU:C	1:A:120:SER:N	2.54	0.58
1:B:87:THR:HG22	1:B:88:ALA:N	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:LYS:CD	1:A:84:LYS:C	2.72	0.58
1:B:195:THR:O	1:B:199:LEU:HB2	2.03	0.58
1:A:294:GLU:OE1	1:A:297:LYS:CE	2.51	0.58
1:B:243:ILE:HG22	1:B:250:THR:HG23	1.85	0.58
1:A:384:ILE:HG22	1:A:385:ARG:N	2.19	0.58
1:A:156:ARG:HG2	1:A:157:ASP:N	2.14	0.57
1:B:390:SER:O	1:B:391:GLY:O	2.22	0.57
1:A:413:GLU:HB3	1:A:416:ILE:HB	1.85	0.57
1:A:401:ILE:HG23	1:A:415:GLU:OE2	2.04	0.57
1:B:430:LEU:O	1:B:431:ALA:HB3	2.04	0.57
1:B:169:LEU:C	1:B:173:ASP:OD1	2.43	0.57
1:A:337:TYR:N	1:A:338:PRO:HD2	2.19	0.56
1:B:183:ARG:HG3	1:B:227:GLU:HG3	1.87	0.56
1:B:262:ASN:ND2	1:B:265:ARG:HH22	2.03	0.56
1:B:285:HIS:HD2	1:B:350:HIS:HB2	1.70	0.56
1:B:320:ALA:HB3	1:B:321:PRO:HD3	1.86	0.56
1:B:337:TYR:HB3	1:B:403:MET:HE3	1.87	0.56
1:A:297:LYS:HE3	1:A:298:GLU:HG2	1.86	0.56
1:A:421:LEU:HD12	1:A:424:LEU:HD13	1.88	0.56
1:A:292:ASP:OD1	1:A:295:SER:HB2	2.05	0.56
1:A:33:VAL:HG23	1:A:35:MET:HG3	1.86	0.56
1:A:90:LEU:HD12	1:A:92:LEU:HG	1.87	0.56
1:A:9:LYS:CD	1:A:9:LYS:N	2.65	0.56
1:B:324:LEU:HD12	1:B:359:LYS:HE2	1.88	0.56
1:A:263:LEU:O	1:A:267:ILE:HG13	2.06	0.55
1:B:139:LYS:O	1:B:145:PHE:O	2.24	0.55
1:B:174:ASP:OD1	1:B:174:ASP:C	2.43	0.55
1:A:272:GLU:C	1:A:273:ASN:HD22	2.08	0.55
1:A:84:LYS:CD	1:A:84:LYS:O	2.54	0.55
1:A:211:ARG:NH2	1:A:211:ARG:CG	2.68	0.55
1:A:352:LEU:HA	1:A:355:LEU:HD21	1.87	0.55
1:A:289:GLU:OE1	1:A:350:HIS:HE1	1.89	0.55
1:B:371:LYS:O	1:B:374:LYS:HD3	2.06	0.55
1:A:253:THR:HB	1:A:254:PRO:HD2	1.89	0.55
1:A:104:ARG:HH11	1:A:128:ARG:HD3	1.72	0.55
1:A:9:LYS:HD3	1:A:9:LYS:H	1.67	0.55
1:B:56:ILE:HG23	1:B:58:ILE:CD1	2.37	0.55
1:A:10:LEU:HD11	1:A:14:ARG:NE	2.22	0.55
1:B:231:LEU:O	1:B:235:ARG:HG3	2.07	0.55
1:A:388:ILE:O	1:A:388:ILE:CG2	2.56	0.54
1:B:63:ASN:C	1:B:63:ASN:OD1	2.46	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:TRP:CH2	1:A:364:LYS:HG2	2.42	0.54
1:A:31:ASP:OD2	1:A:142:PRO:HD2	2.07	0.54
1:A:369:LYS:O	1:A:373:THR:HB	2.07	0.54
1:A:117:VAL:C	1:A:118:GLU:HG2	2.28	0.54
1:A:37:VAL:N	1:A:141:ASN:HD21	1.99	0.54
1:A:167:HIS:HD2	1:A:169:LEU:H	1.56	0.54
1:A:278:ASP:OD1	1:A:281:TYR:HD2	1.91	0.54
1:A:285:HIS:ND1	1:A:350:HIS:HD2	2.06	0.54
1:B:420:VAL:O	1:B:423:SER:CB	2.55	0.54
1:B:33:VAL:HG23	1:B:35:MET:HG3	1.90	0.54
1:B:388:ILE:HG22	1:B:389:THR:H	1.68	0.54
1:A:63:ASN:HB3	1:A:66:GLU:HB2	1.89	0.54
1:A:211:ARG:HG2	1:A:211:ARG:NH2	2.23	0.54
1:A:424:LEU:C	1:A:425:GLU:HG2	2.27	0.54
1:B:391:GLY:O	1:B:392:LYS:CB	2.56	0.54
1:A:120:SER:HA	1:A:122:ILE:HD12	1.90	0.53
1:A:37:VAL:H	1:A:141:ASN:ND2	1.99	0.53
1:B:169:LEU:O	1:B:173:ASP:OD1	2.25	0.53
1:B:8:SER:HB3	1:B:146:GLY:HA3	1.89	0.53
1:A:410:THR:CG2	1:A:411:ARG:N	2.29	0.53
1:B:380:GLY:C	1:B:381:GLU:CG	2.76	0.53
1:A:331:VAL:O	1:A:366:TYR:OH	2.21	0.53
1:B:369:LYS:NZ	1:B:373:THR:HG21	2.23	0.53
1:B:381:GLU:O	1:B:383:LEU:N	2.41	0.53
1:A:90:LEU:HB3	1:A:98:ILE:HG22	1.89	0.53
1:B:381:GLU:C	1:B:383:LEU:N	2.62	0.53
1:A:167:HIS:CD2	1:A:169:LEU:H	2.27	0.53
1:A:289:GLU:OE1	1:A:350:HIS:CE1	2.62	0.53
1:A:382:TYR:O	1:A:383:LEU:C	2.46	0.53
1:A:172:VAL:HG22	1:A:208:TYR:CZ	2.45	0.52
1:A:208:TYR:O	1:A:212:THR:HG22	2.10	0.52
1:B:425:GLU:C	1:B:426:THR:CG2	2.70	0.52
1:B:56:ILE:HG23	1:B:58:ILE:HD12	1.91	0.52
1:B:236:ARG:HD2	1:B:240:PHE:CE2	2.44	0.52
1:B:374:LYS:O	1:B:376:GLU:N	2.42	0.52
1:B:381:GLU:O	1:B:384:ILE:N	2.42	0.52
1:B:431:ALA:O	1:B:432:ALA:C	2.47	0.52
1:A:22:LEU:O	1:A:26:LEU:HG	2.09	0.52
1:A:411:ARG:O	1:A:412:ASP:C	2.48	0.52
1:B:154:GLY:O	1:B:158:LEU:HB2	2.10	0.52
1:B:172:VAL:HG13	1:B:208:TYR:HE1	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:ASN:HD22	1:B:265:ARG:NH1	2.04	0.52
1:B:247:PHE:HB2	1:B:250:THR:HG22	1.92	0.52
1:A:10:LEU:HD11	1:A:14:ARG:CZ	2.39	0.51
1:B:382:TYR:HA	1:B:385:ARG:HB2	1.92	0.51
1:B:185:GLU:OE2	1:B:192:ILE:HG13	2.09	0.51
1:A:373:THR:C	1:A:375:LEU:H	2.14	0.51
1:A:61:GLU:OE1	1:A:119:MET:O	2.29	0.51
1:B:177:ARG:HA	1:B:180:ARG:HB2	1.92	0.51
1:A:63:ASN:OD1	1:A:86:MET:CE	2.58	0.51
1:A:63:ASN:ND2	1:A:65:LEU:HB2	2.25	0.51
1:B:310:ILE:HG23	1:B:311:ASN:N	2.26	0.51
1:B:247:PHE:HB2	1:B:250:THR:CG2	2.40	0.51
1:A:38:TYR:CZ	1:A:122:ILE:HG12	2.46	0.50
1:A:175:PRO:HB2	1:A:212:THR:HG21	1.92	0.50
1:B:418:GLU:O	1:B:421:LEU:HD22	2.11	0.50
1:B:420:VAL:O	1:B:423:SER:N	2.43	0.50
1:B:91:PHE:HA	1:B:97:ARG:HB2	1.94	0.50
1:B:327:LEU:HD11	1:B:363:PHE:HB2	1.93	0.50
1:B:77:GLY:HA3	1:B:91:PHE:O	2.11	0.50
1:A:182:ILE:HG23	1:A:236:ARG:HG2	1.92	0.50
1:A:34:ASN:OD1	1:A:34:ASN:O	2.30	0.50
1:B:306:ARG:NH2	1:B:312:GLU:OE1	2.45	0.50
1:B:388:ILE:HG23	1:B:389:THR:H	1.76	0.50
1:B:38:TYR:CE2	1:B:139:LYS:HG3	2.46	0.50
1:B:56:ILE:CG2	1:B:58:ILE:CD1	2.89	0.50
1:B:63:ASN:O	1:B:63:ASN:OD1	2.30	0.50
1:B:73:ARG:NH1	3:B:443:HOH:O	2.42	0.50
1:A:90:LEU:HB3	1:A:98:ILE:HG23	1.89	0.50
1:A:269:TRP:CZ2	1:A:364:LYS:HG2	2.47	0.50
1:A:272:GLU:CB	1:A:273:ASN:ND2	2.71	0.50
1:B:293:ASP:HA	1:B:314:ARG:HH12	1.77	0.50
1:A:168:THR:HG22	1:A:169:LEU:HG	1.93	0.49
1:A:166:LEU:N	1:A:166:LEU:HD23	2.26	0.49
1:A:191:ARG:HH21	1:A:191:ARG:HG2	1.77	0.49
1:A:415:GLU:HA	1:A:418:GLU:HB2	1.94	0.49
1:B:332:PRO:HB2	1:B:407:ASP:OD1	2.12	0.49
1:A:63:ASN:OD1	1:A:86:MET:HE2	2.12	0.49
1:B:169:LEU:O	1:B:172:VAL:N	2.45	0.49
1:B:374:LYS:O	1:B:402:LEU:HD23	2.12	0.49
1:B:174:ASP:OD1	1:B:174:ASP:O	2.30	0.49
1:B:391:GLY:O	1:B:392:LYS:HD2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:LYS:CD	1:A:144:ASP:HB2	2.38	0.49
1:A:409:ASP:OD1	1:A:409:ASP:C	2.50	0.49
1:B:168:THR:HA	1:B:199:LEU:HD21	1.94	0.49
1:B:296:TRP:CD2	1:B:310:ILE:HB	2.47	0.49
1:B:164:ARG:HB2	1:B:193:GLU:HG2	1.95	0.49
1:A:36:PRO:HG2	1:A:119:MET:HG2	1.94	0.49
1:B:337:TYR:CD2	1:B:403:MET:HE1	2.48	0.49
1:A:396:GLU:O	1:A:400:LYS:HG3	2.13	0.49
1:B:165:VAL:HG22	1:B:196:THR:HG23	1.95	0.49
1:A:104:ARG:O	1:A:105:LEU:C	2.51	0.49
1:A:410:THR:HB	1:A:415:GLU:HB3	1.94	0.49
1:A:172:VAL:HG22	1:A:208:TYR:CE1	2.48	0.48
1:B:203:ALA:O	1:B:208:TYR:HB2	2.13	0.48
1:B:262:ASN:HD22	1:B:265:ARG:HH22	1.59	0.48
2:B:502:ATP:H8	2:B:502:ATP:O5'	1.97	0.48
1:B:126:LEU:HB3	1:B:152:PHE:HE2	1.78	0.48
1:B:93:LYS:N	1:B:93:LYS:CD	2.73	0.48
1:B:62:GLY:O	1:B:63:ASN:CB	2.62	0.48
1:A:231:LEU:HD13	1:A:235:ARG:NH2	2.29	0.48
1:A:310:ILE:O	1:A:310:ILE:HG13	2.14	0.48
1:B:141:ASN:ND2	1:B:141:ASN:H	2.11	0.48
1:B:235:ARG:HH11	1:B:257:ASP:CG	2.17	0.48
1:B:230:PRO:HB2	1:B:264:PHE:CZ	2.49	0.48
1:B:411:ARG:HG3	1:B:413:GLU:H	1.79	0.48
1:B:63:ASN:O	1:B:64:ALA:C	2.52	0.48
1:B:409:ASP:OD1	1:B:409:ASP:O	2.31	0.47
1:B:183:ARG:NH2	1:B:220:GLU:OE1	2.47	0.47
1:B:205:GLU:C	1:B:207:GLY:H	2.16	0.47
1:A:90:LEU:CD1	1:A:90:LEU:C	2.83	0.47
1:B:411:ARG:O	1:B:412:ASP:CB	2.63	0.47
1:B:82:HIS:HB2	1:B:87:THR:HB	1.95	0.47
1:B:285:HIS:CD2	1:B:350:HIS:HB2	2.49	0.47
1:B:234:ILE:HA	1:B:237:MET:HE2	1.97	0.47
1:B:421:LEU:C	1:B:423:SER:N	2.63	0.47
1:A:193:GLU:OE2	1:A:195:THR:HB	2.15	0.47
1:B:188:PHE:HB3	1:B:190:PHE:CE2	2.50	0.47
1:B:381:GLU:C	1:B:383:LEU:H	2.18	0.47
1:B:56:ILE:HG22	1:B:58:ILE:HD12	1.97	0.47
1:A:389:THR:HB	1:A:394:ILE:CD1	2.44	0.47
1:B:382:TYR:HA	1:B:385:ARG:CB	2.44	0.47
1:B:425:GLU:O	1:B:426:THR:CB	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:PRO:O	1:A:92:LEU:HD22	2.14	0.46
1:A:331:VAL:CB	1:A:332:PRO:HD2	2.23	0.46
1:B:15:VAL:HG12	1:B:16:ASP:O	2.16	0.46
1:A:329:GLU:O	1:A:331:VAL:HG13	2.14	0.46
1:B:337:TYR:HD2	1:B:403:MET:HE1	1.81	0.46
1:A:377:LYS:NZ	1:A:377:LYS:HB2	2.31	0.46
1:A:49:LEU:CD2	1:A:158:LEU:HD22	2.42	0.46
1:A:3:ILE:O	1:A:3:ILE:HG22	2.15	0.46
1:B:421:LEU:HD23	1:B:422:ALA:N	2.31	0.46
1:A:329:GLU:O	1:A:330:ARG:C	2.50	0.46
1:A:178:ILE:HD12	1:A:209:LEU:HD13	1.98	0.45
1:B:416:ILE:O	1:B:420:VAL:HG22	2.15	0.45
1:A:117:VAL:HG23	1:A:118:GLU:CG	2.47	0.45
1:A:191:ARG:O	1:A:191:ARG:HG3	2.15	0.45
1:A:28:LYS:HE3	1:B:145:PHE:CG	2.51	0.45
1:B:280:PHE:CZ	1:B:284:LEU:HD11	2.52	0.45
1:A:389:THR:HB	1:A:394:ILE:HD11	1.97	0.45
1:A:56:ILE:CG1	1:A:98:ILE:HG12	2.47	0.45
1:A:36:PRO:HD2	1:A:61:GLU:HB2	1.98	0.45
1:B:129:ARG:HG2	2:B:502:ATP:H1'	1.98	0.45
1:B:169:LEU:O	1:B:170:SER:C	2.54	0.45
1:A:217:LEU:HA	1:A:217:LEU:HD12	1.87	0.45
1:A:394:ILE:O	1:A:397:VAL:HG22	2.16	0.45
1:B:209:LEU:H	1:B:209:LEU:CD1	2.28	0.45
1:B:379:ASN:O	1:B:379:ASN:ND2	2.46	0.45
1:B:413:GLU:O	1:B:416:ILE:CG2	2.62	0.45
1:A:185:GLU:HG2	1:A:185:GLU:O	2.16	0.45
1:A:49:LEU:HB3	1:A:51:ILE:CD1	2.46	0.45
1:A:179:LEU:HA	1:A:179:LEU:HD23	1.76	0.45
1:A:405:LYS:O	1:A:408:GLY:O	2.35	0.45
1:A:285:HIS:HA	1:A:313:ILE:HG12	1.99	0.45
1:B:398:LEU:HA	1:B:398:LEU:HD12	1.84	0.45
1:A:34:ASN:CG	1:A:34:ASN:O	2.54	0.44
1:B:322:ALA:O	1:B:326:MET:HG3	2.17	0.44
1:A:272:GLU:HB2	1:A:273:ASN:HD21	1.74	0.44
1:A:32:GLU:OE2	1:A:70:TYR:OH	2.20	0.44
1:A:394:ILE:O	1:A:398:LEU:HD23	2.18	0.44
1:A:400:LYS:HD3	1:A:419:GLU:OE1	2.17	0.44
1:B:234:ILE:HA	1:B:237:MET:CE	2.47	0.44
1:B:82:HIS:O	1:B:84:LYS:O	2.35	0.44
1:A:231:LEU:HA	1:A:231:LEU:HD23	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:VAL:HG21	1:B:349:CYS:SG	2.58	0.44
1:B:325:GLU:HA	1:B:328:SER:OG	2.17	0.44
1:A:90:LEU:HD12	1:A:92:LEU:CD2	2.35	0.44
1:B:420:VAL:C	1:B:423:SER:H	2.21	0.44
1:B:231:LEU:HD22	1:B:235:ARG:HD2	2.00	0.44
1:B:429:LYS:HE2	1:B:429:LYS:HB2	1.52	0.44
1:A:122:ILE:HD12	1:A:123:LYS:H	1.83	0.44
1:A:250:THR:OG1	1:A:287:PHE:O	2.34	0.44
1:B:369:LYS:O	1:B:373:THR:OG1	2.29	0.44
1:B:296:TRP:CH2	1:B:300:ARG:HB2	2.53	0.44
1:A:270:VAL:O	1:A:271:GLU:C	2.55	0.44
1:A:90:LEU:HD12	1:A:92:LEU:CG	2.47	0.44
1:B:285:HIS:CD2	1:B:350:HIS:HD2	2.35	0.44
1:A:410:THR:HG21	1:A:414:GLU:CB	2.46	0.43
1:B:331:VAL:HG12	1:B:332:PRO:CD	2.43	0.43
1:B:266:ASN:HB3	1:B:352:LEU:HD13	2.01	0.43
1:B:404:LYS:HB2	1:B:415:GLU:OE2	2.18	0.43
1:A:336:VAL:HG11	1:A:366:TYR:CD1	2.53	0.43
1:A:82:HIS:CB	1:A:87:THR:HG22	2.42	0.43
1:B:352:LEU:HA	1:B:355:LEU:HD23	2.00	0.43
1:A:272:GLU:C	1:A:273:ASN:ND2	2.70	0.43
1:A:405:LYS:HE2	1:A:405:LYS:HB3	1.69	0.43
1:B:423:SER:C	1:B:425:GLU:H	2.20	0.43
1:B:93:LYS:HE3	1:B:93:LYS:HB3	1.61	0.43
1:A:63:ASN:ND2	1:A:66:GLU:H	2.15	0.43
1:A:72:LYS:HD2	1:A:72:LYS:HA	1.69	0.43
1:B:20:LEU:HA	1:B:20:LEU:HD23	1.90	0.43
1:B:5:ARG:HG2	1:B:5:ARG:NH1	2.33	0.43
1:A:424:LEU:C	1:A:424:LEU:HD23	2.39	0.43
1:B:430:LEU:O	1:B:431:ALA:CB	2.67	0.43
1:B:9:LYS:HD3	1:B:9:LYS:H	1.83	0.43
1:A:40:VAL:HG21	1:A:59:VAL:HG23	2.00	0.43
1:B:243:ILE:HG22	1:B:250:THR:CG2	2.48	0.43
1:B:92:LEU:HA	1:B:93:LYS:NZ	2.33	0.43
1:B:337:TYR:CZ	1:B:341:LYS:HG3	2.54	0.43
1:B:378:ILE:HG21	1:B:411:ARG:NH2	2.30	0.43
1:B:90:LEU:HD12	1:B:92:LEU:HD21	2.01	0.43
1:A:262:ASN:HD22	1:A:265:ARG:NH2	2.06	0.43
1:A:82:HIS:CG	1:A:87:THR:HG22	2.53	0.43
1:A:352:LEU:HA	1:A:355:LEU:CD2	2.48	0.43
1:A:336:VAL:HG11	1:A:366:TYR:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:GLY:HA3	1:A:140:LEU:O	2.18	0.42
1:B:175:PRO:CD	1:B:212:THR:HG21	2.45	0.42
1:A:105:LEU:HD12	1:A:105:LEU:H	1.84	0.42
1:A:40:VAL:CG2	1:A:59:VAL:HG23	2.48	0.42
1:B:285:HIS:CD2	1:B:350:HIS:CD2	3.08	0.42
1:A:302:ARG:HG2	1:A:303:TYR:CD1	2.54	0.42
1:A:308:ASN:O	1:A:312:GLU:HG3	2.19	0.42
1:A:343:VAL:HB	1:A:347:THR:HB	2.01	0.42
1:B:421:LEU:C	1:B:423:SER:H	2.23	0.42
1:A:300:ARG:O	1:A:303:TYR:O	2.37	0.42
1:B:289:GLU:OE1	1:B:350:HIS:HE1	2.02	0.42
1:B:54:LEU:O	1:B:55:ASP:C	2.58	0.42
1:A:266:ASN:HB3	1:A:352:LEU:HD13	2.01	0.42
1:B:262:ASN:HD22	1:B:265:ARG:NH2	2.18	0.42
1:B:360:GLU:HG3	1:B:361:GLY:N	2.35	0.42
1:B:39:VAL:HG13	1:B:138:ILE:HB	2.00	0.42
1:B:87:THR:CG2	1:B:88:ALA:N	2.83	0.42
1:A:341:LYS:H	1:A:341:LYS:HG2	1.61	0.42
1:A:120:SER:HA	1:A:122:ILE:HD11	2.01	0.42
1:B:357:GLY:O	1:B:360:GLU:HG2	2.20	0.42
1:A:163:ILE:O	1:A:163:ILE:HG22	2.20	0.42
1:A:269:TRP:CD2	1:A:364:LYS:HE2	2.55	0.42
1:B:327:LEU:CD1	1:B:363:PHE:HB2	2.50	0.42
1:A:369:LYS:HE3	1:A:406:LEU:O	2.19	0.42
1:B:84:LYS:O	1:B:85:PHE:HB2	2.20	0.41
1:A:253:THR:HB	1:A:254:PRO:CD	2.49	0.41
1:B:187:ARG:HD2	1:B:188:PHE:CZ	2.55	0.41
1:B:271:GLU:HA	1:B:275:GLY:O	2.19	0.41
1:B:285:HIS:HD2	1:B:350:HIS:CB	2.32	0.41
1:B:388:ILE:C	1:B:389:THR:HG23	2.37	0.41
1:B:75:LEU:N	1:B:75:LEU:HD12	2.34	0.41
1:A:325:GLU:O	1:A:328:SER:O	2.39	0.41
1:B:178:ILE:O	1:B:182:ILE:HG13	2.20	0.41
1:B:91:PHE:CE1	1:B:97:ARG:HD2	2.55	0.41
1:A:195:THR:HG22	1:A:199:LEU:HD22	2.03	0.41
1:B:169:LEU:O	1:B:171:PHE:N	2.53	0.41
1:B:26:LEU:HD23	1:B:26:LEU:HA	1.93	0.41
1:A:36:PRO:CG	1:A:119:MET:HG2	2.50	0.41
1:A:335:PHE:HD2	1:A:335:PHE:O	2.04	0.41
1:A:90:LEU:CD1	1:A:92:LEU:CG	2.94	0.41
2:B:502:ATP:O2B	2:B:502:ATP:O3G	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:ILE:HD12	1:A:209:LEU:CD1	2.50	0.41
1:A:143:LYS:HE3	1:A:143:LYS:HB3	1.69	0.40
1:A:419:GLU:HA	1:A:422:ALA:HB3	2.04	0.40
1:B:209:LEU:HD12	1:B:209:LEU:N	2.28	0.40
1:B:205:GLU:C	1:B:207:GLY:N	2.74	0.40
1:B:411:ARG:O	1:B:412:ASP:HB2	2.20	0.40
1:B:424:LEU:HA	1:B:424:LEU:HD12	1.68	0.40
1:B:355:LEU:HD22	1:B:355:LEU:N	2.36	0.40
1:A:302:ARG:HG2	1:A:303:TYR:CZ	2.56	0.40
1:B:343:VAL:HB	1:B:347:THR:CB	2.52	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	411/441 (93%)	385 (94%)	23 (6%)	3 (1%)	22	50
1	B	416/441 (94%)	378 (91%)	27 (6%)	11 (3%)	5	17
All	All	827/882 (94%)	763 (92%)	50 (6%)	14 (2%)	9	27

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	ASP
1	B	63	ASN
1	B	389	THR
1	B	391	GLY
1	B	426	THR
1	B	429	LYS
1	B	430	LEU
1	B	431	ALA

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Mol	Chain	Res	Type
1	A	118	GLU
1	B	412	ASP
1	A	86	MET
1	B	428	GLY
1	B	329	GLU
1	B	392	LYS

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	374/395 (95%)	320 (86%)	54 (14%)	3	8
1	B	376/395 (95%)	322 (86%)	54 (14%)	3	8
All	All	750/790 (95%)	642 (86%)	108 (14%)	3	8

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	12	VAL
1	A	20	LEU
1	A	28	LYS
1	A	56	ILE
1	A	58	ILE
1	A	63	ASN
1	A	69	GLU
1	A	72	LYS
1	A	75	LEU
1	A	84	LYS
1	A	85	PHE
1	A	87	THR
1	A	98	ILE
1	A	104	ARG
1	A	118	GLU
1	A	119	MET

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Mol	Chain	Res	Type
1	A	122	ILE
1	A	126	LEU
1	A	156	ARG
1	A	163	ILE
1	A	168	THR
1	A	176	THR
1	A	199	LEU
1	A	205	GLU
1	A	211	ARG
1	A	212	THR
1	A	217	LEU
1	A	221	LEU
1	A	223	LYS
1	A	231	LEU
1	A	241	ASP
1	A	263	LEU
1	A	273	ASN
1	A	297	LYS
1	A	302	ARG
1	A	306	ARG
1	A	323	LEU
1	A	324	LEU
1	A	335	PHE
1	A	336	VAL
1	A	340	VAL
1	A	350	HIS
1	A	355	LEU
1	A	367	LEU
1	A	374	LYS
1	A	383	LEU
1	A	384	ILE
1	A	389	THR
1	A	399	GLU
1	A	402	LEU
1	A	407	ASP
1	A	413	GLU
1	A	414	GLU
1	B	9	LYS
1	B	12	VAL
1	B	22	LEU
1	B	28	LYS
1	B	83	ASP

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Mol	Chain	Res	Type
1	B	90	LEU
1	B	93	LYS
1	B	97	ARG
1	B	104	ARG
1	B	118	GLU
1	B	122	ILE
1	B	126	LEU
1	B	148	LEU
1	B	158	LEU
1	B	159	LYS
1	B	189	ASP
1	B	210	GLU
1	B	211	ARG
1	B	213	THR
1	B	223	LYS
1	B	231	LEU
1	B	258	GLU
1	B	263	LEU
1	B	271	GLU
1	B	279	ARG
1	B	285	HIS
1	B	288	LEU
1	B	323	LEU
1	B	329	GLU
1	B	330	ARG
1	B	335	PHE
1	B	370	ILE
1	B	372	ASN
1	B	374	LYS
1	B	375	LEU
1	B	377	LYS
1	B	378	ILE
1	B	379	ASN
1	B	383	LEU
1	B	384	ILE
1	B	385	ARG
1	B	397	VAL
1	B	398	LEU
1	B	403	MET
1	B	407	ASP
1	B	411	ARG
1	B	414	GLU

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Mol	Chain	Res	Type
1	B	417	LEU
1	B	421	LEU
1	B	423	SER
1	B	424	LEU
1	B	426	THR
1	B	429	LYS
1	B	430	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	82	HIS
1	A	141	ASN
1	A	167	HIS
1	A	262	ASN
1	A	273	ASN
1	A	345	ASN
1	A	350	HIS
1	B	2	GLN
1	B	141	ASN
1	B	219	GLN
1	B	229	ASN
1	B	245	HIS
1	B	262	ASN
1	B	285	HIS
1	B	345	ASN
1	B	350	HIS
1	B	372	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	ATP	B	502	-	26,33,33	0.96	1 (3%)	31,52,52	1.33	4 (12%)
2	ATP	A	501	-	26,33,33	1.01	2 (7%)	31,52,52	1.64	6 (19%)

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	ATP	B	502	-	-	5/18/38/38	0/3/3/3
2	ATP	A	501	-	-	1/18/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ATP	C5-C4	2.72	1.48	1.40
2	B	502	ATP	C5-C4	2.55	1.47	1.40
2	A	501	ATP	C2'-C1'	-2.02	1.50	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ATP	C4-C5-N7	-3.66	105.59	109.40
2	A	501	ATP	PB-O3B-PG	-3.61	120.43	132.83
2	B	502	ATP	PB-O3B-PG	-2.93	122.76	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	ATP	N3-C2-N1	-2.85	124.23	128.68
2	A	501	ATP	N3-C2-N1	-2.82	124.28	128.68
2	A	501	ATP	PA-O3A-PB	-2.72	123.49	132.83
2	B	502	ATP	C4-C5-N7	-2.61	106.68	109.40
2	B	502	ATP	PA-O3A-PB	-2.51	124.20	132.83
2	A	501	ATP	O2G-PG-O1G	2.04	118.68	110.68
2	A	501	ATP	O2A-PA-O5'	2.01	117.08	107.75

There are no chirality outliers.

All (6) torsion outliers are listed below:

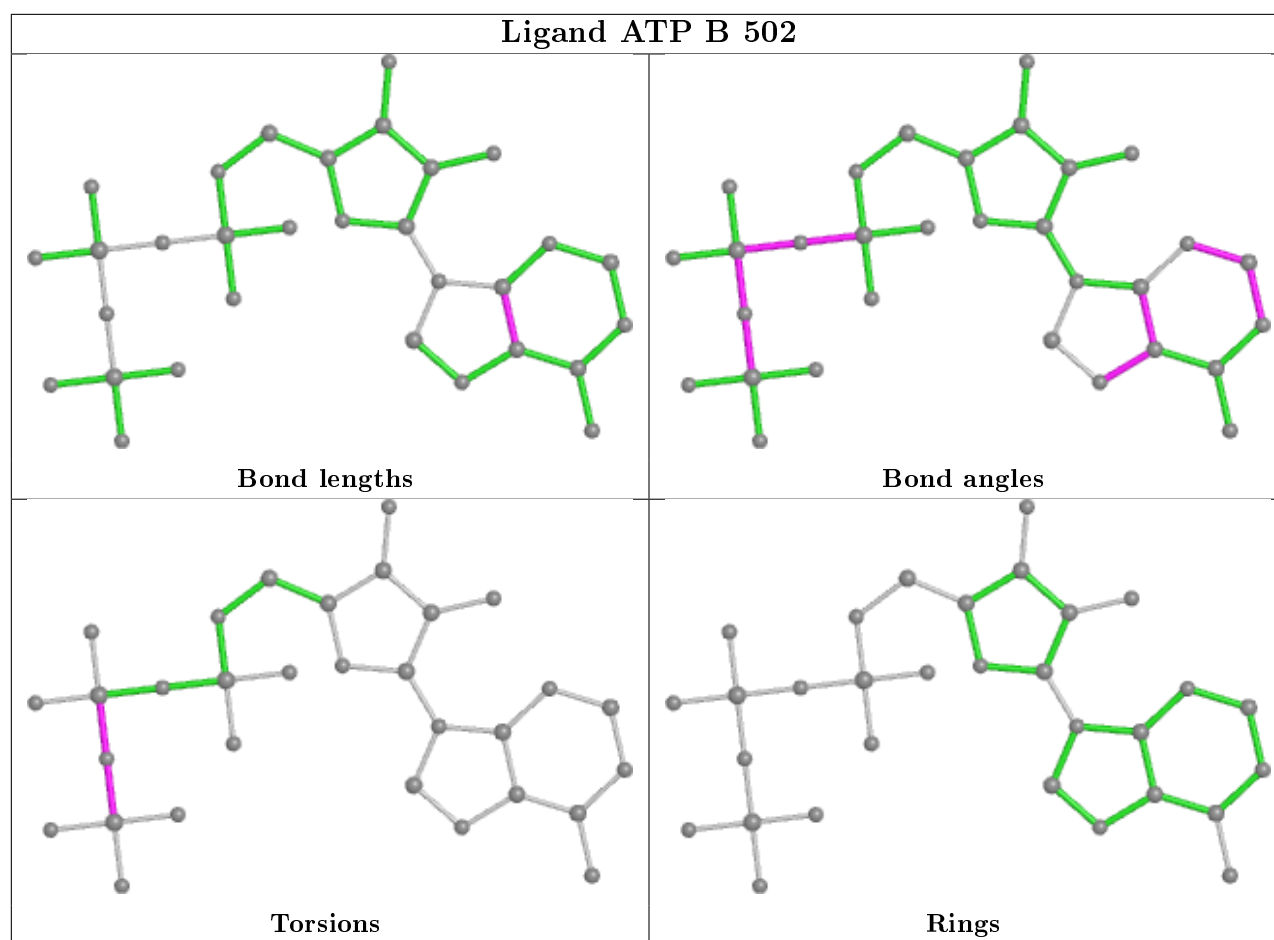
Mol	Chain	Res	Type	Atoms
2	B	502	ATP	PB-O3B-PG-O2G
2	B	502	ATP	PB-O3B-PG-O3G
2	B	502	ATP	PG-O3B-PB-O1B
2	B	502	ATP	PG-O3B-PB-O2B
2	A	501	ATP	O4'-C4'-C5'-O5'
2	B	502	ATP	PB-O3B-PG-O1G

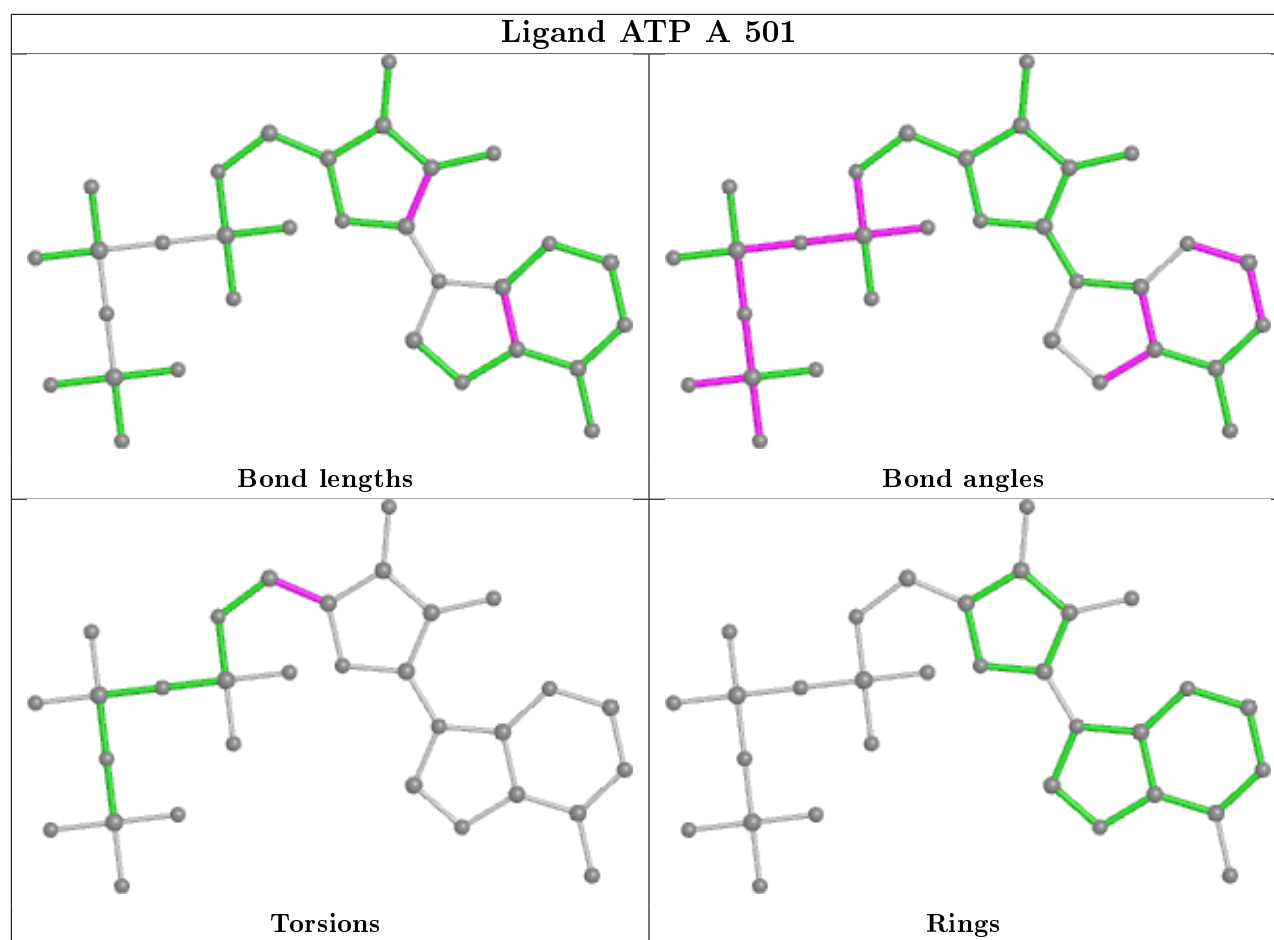
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	ATP	3	0
2	A	501	ATP	1	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	415/441 (94%)	-0.12	12 (2%) 51 47	30, 59, 154, 192	0
1	B	420/441 (95%)	-0.05	9 (2%) 63 60	24, 95, 142, 164	0
All	All	835/882 (94%)	-0.09	21 (2%) 57 54	24, 76, 148, 192	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	432	ALA	5.2
1	A	117	VAL	4.4
1	A	374	LYS	3.8
1	A	378	ILE	3.8
1	B	105	LEU	3.6
1	A	410	THR	3.5
1	A	416	ILE	3.0
1	A	415	GLU	2.9
1	A	375	LEU	2.9
1	A	411	ARG	2.5
1	B	362	LEU	2.4
1	B	119	MET	2.3
1	B	385	ARG	2.3
1	A	381	GLU	2.3
1	B	85	PHE	2.3
1	B	351	PHE	2.3
1	B	414	GLU	2.2
1	A	384	ILE	2.2
1	A	420	VAL	2.0
1	B	415	GLU	2.0
1	A	95	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

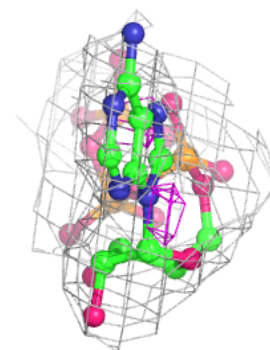
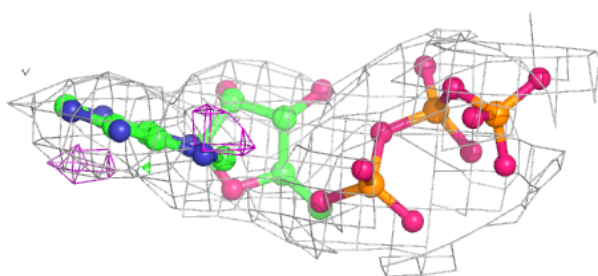
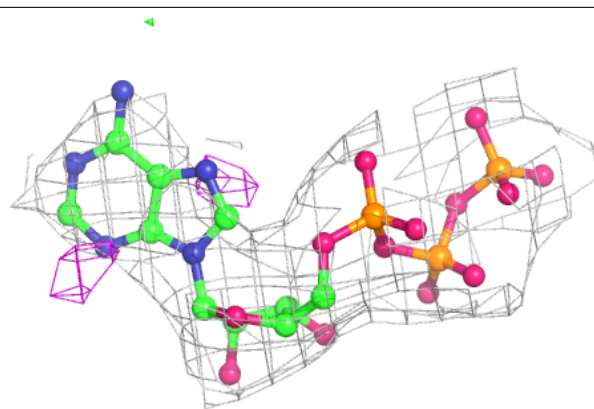
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ATP	B	502	31/31	0.92	0.20	67,99,123,134	0
2	ATP	A	501	31/31	0.97	0.15	26,43,62,102	0

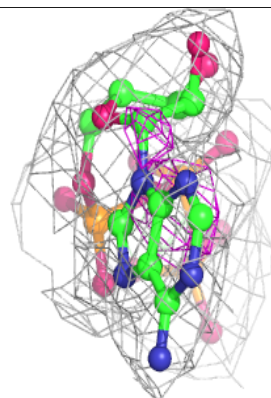
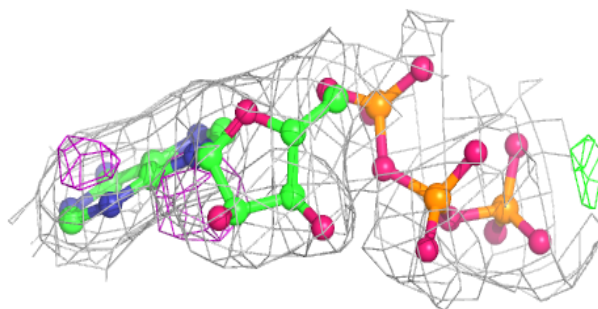
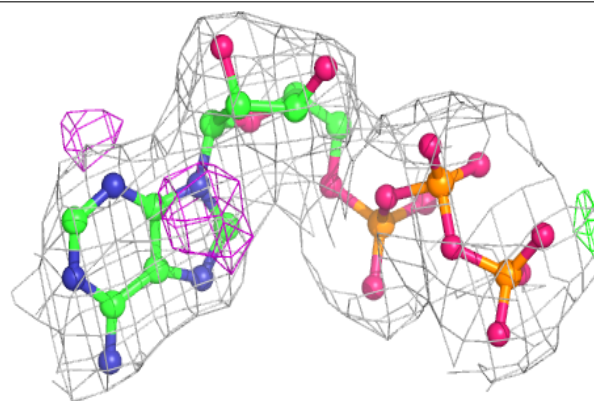
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP B 502:

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

**Electron density around ATP A 501:**

$2mF_o-DF_c$ (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.