



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:02 AM UTC

PDB ID : 3AQN / pdb_00003aqn
Title : Complex structure of bacterial protein (apo form II)
Authors : Toh, Y.; Takeshita, D.; Tomita , K.
Deposited on : 2010-11-09
Resolution : 3.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

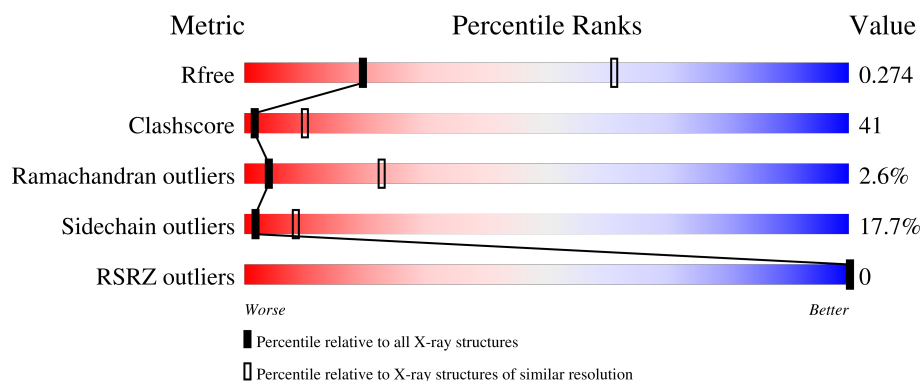
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

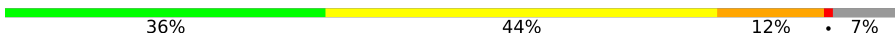
The reported resolution of this entry is 3.30 Å.

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))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	415	
1	B	415	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6398 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 Poly(A) polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			3164	2022	568	560	14			
1	B	388	Total	C	N	O	S	0	0	0
			3164	2022	568	560	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	HIS	ARG	engineered mutation	UNP C9QS13
B	234	HIS	ARG	engineered mutation	UNP C9QS13

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: 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		

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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

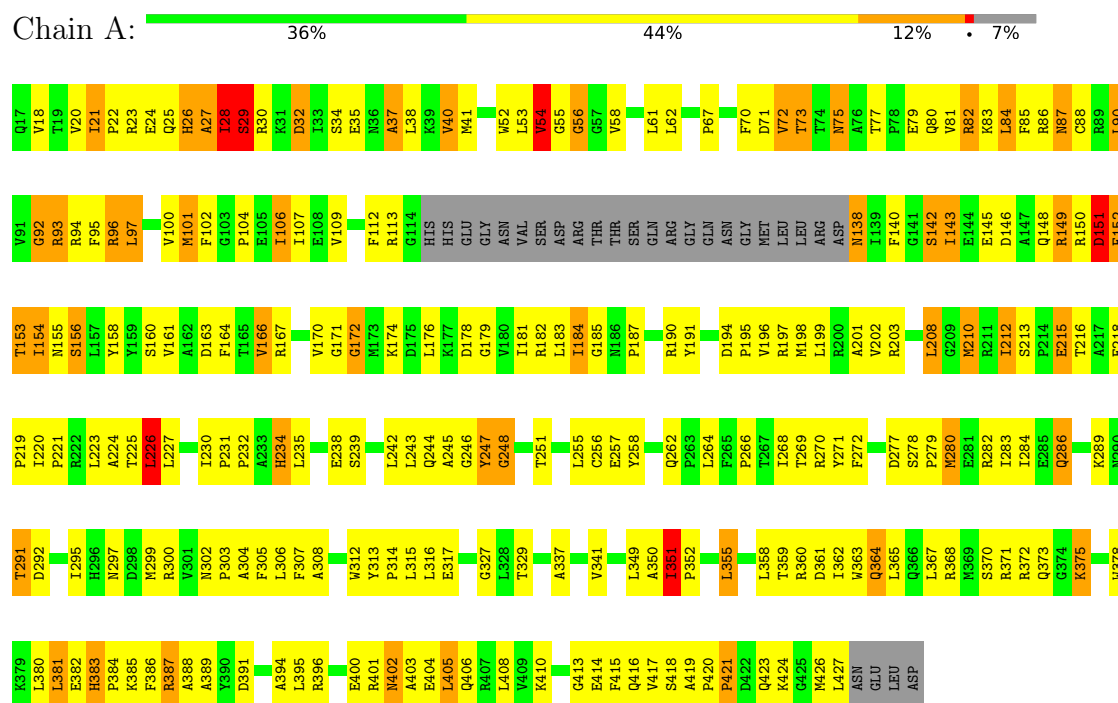
- Molecule 4 is water.

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

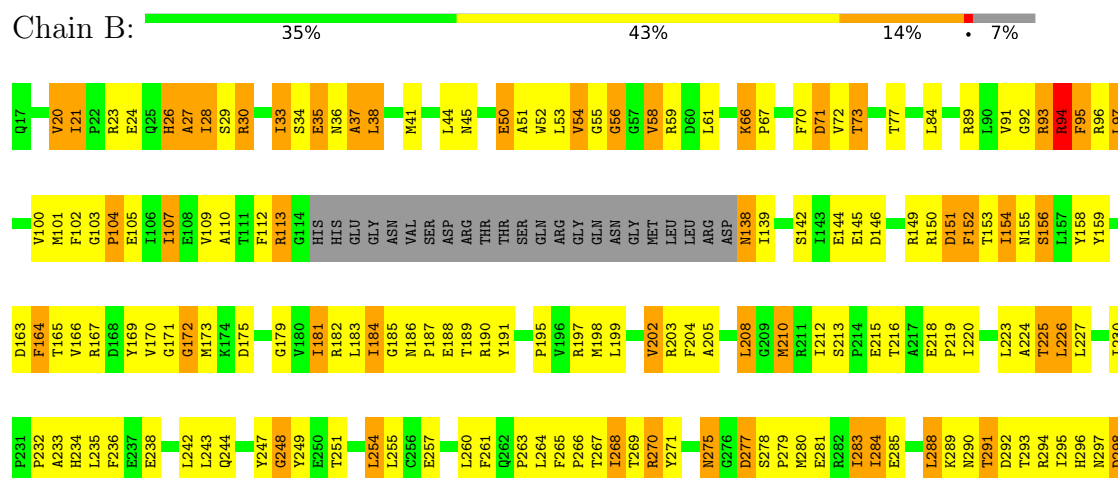
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Poly(A) polymerase



• Molecule 1: Poly(A) polymerase



M299	R368	X369	S370	R371	R372	Q373	G374		W378	K379	L380	L381	E382	H383	P384	K385	F386	R387		D391	L392	L393	A394	L395	R396		N402	A403	E404	L405	Q406	R407	L408	V409		E414	F415	Q416	V417	S418	A419	P420		Q423		L427	ASN	GLU	LEU	ASP																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																					

4 Data and refinement statistics

Property	Value	Source
Space group	P 4 ₂ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	132.40Å 132.40Å 176.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.61 – 3.30 29.61 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.61-3.30) 98.9 (29.61-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 3.19Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.251 , 0.275 0.251 , 0.274	Depositor DCC
R_{free} test set	1364 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	90.5	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 67.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6398	wwPDB-VP
Average B, all atoms (Å ²)	128.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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, MG

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	0.50	0/3234	0.98	19/4371 (0.4%)
1	B	0.49	0/3234	0.97	19/4371 (0.4%)
All	All	0.49	0/6468	0.97	38/8742 (0.4%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	ALA	N-CA-C	-10.82	94.07	110.30
1	A	21	ILE	CA-C-N	10.14	130.09	119.85
1	A	21	ILE	C-N-CA	10.14	130.09	119.85
1	A	151	ASP	N-CA-C	9.45	121.18	111.07
1	A	92	GLY	N-CA-C	7.74	125.31	112.85
1	B	291	THR	N-CA-C	-7.58	102.05	111.75
1	B	406	GLN	N-CA-C	-7.25	102.47	111.75
1	A	153	THR	N-CA-C	7.21	118.78	111.07
1	B	54	VAL	N-CA-C	7.15	118.78	108.48
1	B	95	PHE	N-CA-C	-7.01	98.31	109.59
1	B	383	HIS	CA-C-N	6.92	126.89	119.28
1	B	383	HIS	C-N-CA	6.92	126.89	119.28
1	A	37	ALA	N-CA-C	-6.69	104.99	113.02
1	B	94	ARG	CB-CA-C	6.55	123.46	110.42
1	B	66	LYS	CA-C-N	6.43	126.45	119.90
1	B	66	LYS	C-N-CA	6.43	126.45	119.90
1	B	420	PRO	N-CA-C	6.31	118.40	110.70
1	A	383	HIS	CA-C-N	6.25	125.94	119.56
1	A	383	HIS	C-N-CA	6.25	125.94	119.56
1	B	91	VAL	N-CA-C	6.23	116.68	107.15
1	A	351	ILE	CA-C-N	6.23	126.59	119.93
1	A	351	ILE	C-N-CA	6.23	126.59	119.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	181	ILE	CB-CA-C	-5.92	105.52	111.80
1	B	351	ILE	CA-C-N	5.91	125.84	119.76
1	B	351	ILE	C-N-CA	5.91	125.84	119.76
1	A	104	PRO	N-CA-C	-5.89	106.44	114.80
1	B	94	ARG	N-CA-C	-5.73	98.60	110.80
1	A	54	VAL	N-CA-C	5.67	116.91	108.46
1	B	37	ALA	N-CA-C	-5.51	106.52	113.18
1	A	247	TYR	N-CA-C	-5.50	106.42	114.39
1	A	142	SER	N-CA-C	5.49	118.19	110.23
1	A	97	LEU	N-CA-C	5.43	118.68	111.30
1	B	95	PHE	N-CA-CB	5.35	119.30	110.69
1	B	27	ALA	N-CA-C	-5.32	102.42	110.24
1	B	346	CYS	N-CA-C	-5.30	106.88	113.19
1	A	96	ARG	CB-CA-C	-5.27	110.52	116.63
1	A	245	ALA	N-CA-C	-5.20	106.25	112.59
1	A	152	PHE	N-CA-C	5.07	117.56	109.81

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	3164	0	3197	288	0
1	B	3164	0	3197	249	0
2	A	31	0	12	4	0
2	B	31	0	12	7	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
All	All	6398	0	6418	527	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 41.

All (527) 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:28:ILE:O	1:A:29:SER:C	1.85	1.19
1:A:191:TYR:CD1	1:A:198:MET:HG3	1.88	1.08
1:B:374:GLY:HA2	1:B:427:LEU:HD22	1.37	1.06
1:B:93:ARG:H	1:B:93:ARG:HD3	1.15	1.06
1:A:184:ILE:HG22	1:A:190:ARG:HH12	1.18	1.05
1:A:26:HIS:HD2	1:A:27:ALA:O	1.38	1.04
1:A:28:ILE:CG2	1:A:30:ARG:HG2	1.87	1.03
1:A:96:ARG:HH12	1:A:113:ARG:HD3	1.11	1.03
1:A:95:PHE:O	1:A:96:ARG:CG	2.08	1.02
1:A:28:ILE:HG23	1:A:30:ARG:HG2	1.03	1.01
1:B:183:LEU:HD23	1:B:184:ILE:H	1.24	1.01
1:A:28:ILE:HG23	1:A:30:ARG:CG	1.93	0.99
1:A:191:TYR:HD1	1:A:198:MET:HG3	1.26	0.95
1:A:95:PHE:O	1:A:96:ARG:HG2	1.66	0.95
1:A:21:ILE:HG23	1:A:166:VAL:HG13	1.46	0.94
1:A:153:THR:HG22	1:A:170:VAL:CG1	1.97	0.94
1:A:28:ILE:O	1:A:28:ILE:HD13	1.66	0.93
1:A:154:ILE:HG13	1:A:155:ASN:ND2	1.84	0.93
1:A:96:ARG:NH1	1:A:113:ARG:HD3	1.83	0.92
1:A:79:GLU:HA	1:A:82:ARG:HE	1.31	0.92
1:A:184:ILE:HG22	1:A:190:ARG:NH1	1.85	0.92
1:A:183:LEU:HD23	1:A:184:ILE:H	1.34	0.92
1:B:202:VAL:HG12	1:B:254:LEU:HB3	1.52	0.91
1:A:77:THR:H	1:A:80:GLN:HE21	1.14	0.91
1:A:153:THR:HG22	1:A:170:VAL:HG11	1.54	0.90
1:A:308:ALA:HB2	1:A:359:THR:HG23	1.55	0.88
1:A:212:ILE:H	1:A:212:ILE:HD12	1.35	0.87
1:A:244:GLN:HE22	1:A:303:PRO:HA	1.41	0.86
1:A:34:SER:HB3	1:A:38:LEU:HD21	1.57	0.86
1:B:183:LEU:HB2	1:B:216:THR:HG23	1.57	0.86
1:B:35:GLU:HA	1:B:35:GLU:OE1	1.76	0.86
1:A:26:HIS:CD2	1:A:27:ALA:O	2.28	0.86
1:A:94:ARG:HB3	1:A:97:LEU:HD23	1.58	0.86
1:A:154:ILE:HG13	1:A:155:ASN:HD22	1.40	0.84
1:B:280:MET:HE2	1:B:314:PRO:HD3	1.59	0.84
1:A:96:ARG:HH22	1:A:113:ARG:HB3	1.42	0.83
1:A:208:LEU:HD13	1:A:210:MET:HE2	1.59	0.82
1:B:170:VAL:HG12	1:B:171:GLY:H	1.45	0.82
1:A:304:ALA:HB2	1:A:355:LEU:HD12	1.61	0.81
1:B:308:ALA:HB2	1:B:359:THR:HG23	1.63	0.81
1:B:352:PRO:HG2	1:B:355:LEU:HD23	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:ARG:HD3	1:B:93:ARG:N	1.93	0.80
1:A:26:HIS:O	1:A:27:ALA:C	2.25	0.80
1:A:95:PHE:O	1:A:96:ARG:HG3	1.81	0.80
1:A:28:ILE:CG2	1:A:30:ARG:HE	1.95	0.79
1:A:92:GLY:C	1:A:93:ARG:HD2	2.08	0.79
1:A:191:TYR:CE1	1:A:198:MET:HG3	2.19	0.77
1:A:86:ARG:HG3	1:A:87:ASN:HD22	1.48	0.77
1:A:95:PHE:C	1:A:96:ARG:HG2	2.09	0.77
1:B:203:ARG:HB2	1:B:242:LEU:HD23	1.66	0.77
1:A:183:LEU:HD23	1:A:184:ILE:N	1.99	0.77
1:B:270:ARG:HD2	1:B:271:TYR:CE2	2.20	0.77
1:B:232:PRO:HB2	1:B:350:ALA:HB2	1.67	0.76
1:A:28:ILE:O	1:A:29:SER:O	2.03	0.76
1:A:402:ASN:HD22	1:A:403:ALA:N	1.83	0.76
1:A:270:ARG:HD2	1:A:271:TYR:CE2	2.21	0.76
1:B:302:ASN:HD22	1:B:303:PRO:HD2	1.51	0.76
1:B:202:VAL:CG1	1:B:254:LEU:HB3	2.16	0.74
1:A:170:VAL:HG12	1:A:171:GLY:H	1.53	0.74
1:A:94:ARG:CZ	1:A:97:LEU:CD2	2.66	0.73
1:A:153:THR:HG22	1:A:170:VAL:HG12	1.69	0.73
1:A:297:ASN:HD22	1:B:378:TRP:NE1	1.87	0.72
1:A:257:GLU:HG2	1:A:257:GLU:O	1.89	0.72
1:A:218:GLU:HB3	1:A:219:PRO:HD3	1.71	0.72
1:A:151:ASP:OD1	2:A:501:ATP:N3	2.22	0.72
1:B:142:SER:HB3	1:B:145:GLU:HG3	1.72	0.71
1:A:28:ILE:O	1:A:30:ARG:N	2.22	0.71
1:B:170:VAL:HG12	1:B:171:GLY:N	2.05	0.71
1:B:151:ASP:OD1	2:B:502:ATP:C2	2.44	0.71
1:B:244:GLN:HE22	1:B:303:PRO:HA	1.56	0.71
1:A:151:ASP:OD1	2:A:501:ATP:C2	2.44	0.71
1:A:270:ARG:HD2	1:A:271:TYR:HE2	1.56	0.70
1:A:72:VAL:HG22	1:A:109:VAL:HG22	1.72	0.70
1:B:26:HIS:O	1:B:27:ALA:C	2.34	0.70
1:B:34:SER:HB3	1:B:38:LEU:HD21	1.75	0.69
1:B:342:LEU:O	1:B:346:CYS:HB2	1.92	0.69
1:A:27:ALA:O	1:A:29:SER:N	2.22	0.69
1:B:204:PHE:O	1:B:208:LEU:HD12	1.92	0.69
1:A:378:TRP:HB3	1:B:299:MET:SD	2.32	0.69
1:B:93:ARG:HG2	1:B:94:ARG:H	1.56	0.69
1:B:27:ALA:O	1:B:29:SER:N	2.26	0.69
1:B:95:PHE:O	1:B:96:ARG:HG2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:ASP:OD1	2:B:502:ATP:N3	2.26	0.68
1:A:96:ARG:NH2	1:A:113:ARG:HB3	2.06	0.68
1:A:94:ARG:CB	1:A:97:LEU:HD23	2.22	0.68
1:B:28:ILE:O	1:B:29:SER:C	2.36	0.68
1:B:324:GLN:HG2	1:B:325:GLU:N	2.07	0.68
1:A:154:ILE:CG1	1:A:155:ASN:N	2.54	0.68
1:A:154:ILE:CG1	1:A:155:ASN:ND2	2.56	0.68
1:A:113:ARG:HG2	1:A:146:ASP:OD1	1.92	0.68
1:B:218:GLU:HB3	1:B:219:PRO:HD3	1.76	0.68
1:B:95:PHE:O	1:B:96:ARG:CG	2.42	0.68
1:A:149:ARG:HH11	1:A:149:ARG:HB3	1.59	0.67
1:B:35:GLU:OE1	1:B:35:GLU:CA	2.41	0.67
1:B:271:TYR:HE1	1:B:317:GLU:HG2	1.58	0.67
1:A:154:ILE:HG12	1:A:155:ASN:H	1.59	0.67
1:A:28:ILE:HG21	1:A:30:ARG:NE	2.10	0.67
1:A:84:LEU:C	1:A:85:PHE:CD1	2.73	0.67
1:A:154:ILE:HG12	1:A:155:ASN:N	2.10	0.67
1:A:28:ILE:HG21	1:A:30:ARG:HE	1.59	0.67
1:A:94:ARG:NE	1:A:97:LEU:HD23	2.10	0.66
1:A:94:ARG:CG	1:A:97:LEU:HD23	2.26	0.66
1:B:321:LYS:HE2	1:B:325:GLU:OE1	1.95	0.66
1:B:93:ARG:H	1:B:93:ARG:CD	1.99	0.66
1:B:290:ASN:O	1:B:291:THR:C	2.40	0.65
1:A:199:LEU:O	1:A:202:VAL:HG12	1.97	0.65
1:A:420:PRO:O	1:A:424:LYS:HB2	1.96	0.64
1:A:84:LEU:C	1:A:85:PHE:HD1	2.05	0.64
1:A:184:ILE:CG2	1:A:185:GLY:N	2.60	0.64
1:A:358:LEU:HD23	1:A:385:LYS:HE2	1.79	0.64
1:B:191:TYR:CD1	1:B:198:MET:HG3	2.33	0.64
1:B:271:TYR:HB2	1:B:280:MET:HE1	1.78	0.64
1:A:187:PRO:O	1:A:191:TYR:HD2	1.80	0.64
1:B:183:LEU:HB2	1:B:216:THR:CG2	2.28	0.64
1:B:195:PRO:HD2	1:B:230:ILE:HD11	1.80	0.64
1:A:391:ASP:O	1:A:395:LEU:HD13	1.97	0.63
1:B:153:THR:HG22	1:B:170:VAL:CG1	2.27	0.63
1:A:219:PRO:O	1:A:223:LEU:HD12	1.99	0.63
1:A:232:PRO:HB2	1:A:350:ALA:HB2	1.79	0.63
1:A:151:ASP:N	2:A:501:ATP:O2'	2.32	0.63
1:A:300:ARG:HG2	1:B:382:GLU:OE2	1.97	0.63
1:A:271:TYR:HE1	1:A:317:GLU:HG2	1.63	0.63
1:B:144:GLU:HA	1:B:169:TYR:CD2	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:SER:HB3	1:A:163:ASP:OD1	1.98	0.63
1:B:26:HIS:NE2	1:B:29:SER:HB3	2.14	0.63
1:B:243:LEU:HA	1:B:248:GLY:HA2	1.81	0.63
2:A:501:ATP:H5'1	2:A:501:ATP:O1B	1.99	0.63
1:B:44:LEU:HD23	1:B:84:LEU:HD23	1.81	0.62
2:B:502:ATP:H5'1	2:B:502:ATP:O1B	1.99	0.62
1:A:37:ALA:O	1:A:40:VAL:HG13	1.99	0.62
1:B:280:MET:HE2	1:B:314:PRO:CD	2.29	0.62
1:A:24:GLU:HG2	1:A:25:GLN:N	2.14	0.62
1:A:299:MET:SD	1:B:378:TRP:HE3	2.22	0.62
1:A:34:SER:CB	1:A:38:LEU:HD21	2.28	0.62
1:B:33:ILE:HG22	1:B:33:ILE:O	1.99	0.62
1:B:93:ARG:CG	1:B:94:ARG:H	2.11	0.62
1:B:183:LEU:HD23	1:B:184:ILE:N	2.06	0.62
1:A:299:MET:SD	1:B:378:TRP:CE3	2.93	0.62
1:A:153:THR:HB	1:A:172:GLY:H	1.65	0.62
1:B:72:VAL:CG2	1:B:109:VAL:HG22	2.30	0.62
1:B:102:PHE:HB2	1:B:105:GLU:O	1.99	0.61
1:B:203:ARG:HB2	1:B:242:LEU:CD2	2.29	0.61
1:A:28:ILE:CG2	1:A:30:ARG:NE	2.64	0.61
1:A:28:ILE:O	1:A:28:ILE:CD1	2.44	0.61
1:B:153:THR:HG21	1:B:175:ASP:OD2	1.99	0.61
1:A:183:LEU:HB2	1:A:216:THR:HG23	1.82	0.61
1:A:75:ASN:C	1:A:75:ASN:HD22	2.09	0.60
1:B:244:GLN:NE2	1:B:303:PRO:HA	2.17	0.60
1:B:358:LEU:HD23	1:B:385:LYS:HE3	1.83	0.60
1:A:95:PHE:C	1:A:96:ARG:CG	2.68	0.60
1:B:372:ARG:HE	1:B:407:ARG:HH21	1.49	0.60
1:B:381:LEU:HD22	1:B:415:PHE:CE2	2.37	0.60
1:B:352:PRO:CG	1:B:355:LEU:HD23	2.32	0.60
1:A:291:THR:O	1:A:295:ILE:HG12	2.02	0.60
1:B:30:ARG:HE	1:B:33:ILE:HG13	1.67	0.60
1:B:257:GLU:O	1:B:257:GLU:HG2	1.99	0.60
1:A:92:GLY:C	1:A:93:ARG:CD	2.75	0.60
1:B:164:PHE:CD1	1:B:164:PHE:N	2.70	0.59
1:B:37:ALA:CB	1:B:107:ILE:HG12	2.32	0.59
1:A:67:PRO:HG2	1:A:70:PHE:CZ	2.37	0.59
1:A:195:PRO:HB2	1:A:230:ILE:HD11	1.84	0.59
1:A:195:PRO:O	1:A:198:MET:HE3	2.03	0.59
1:A:77:THR:O	1:A:81:VAL:HG23	2.03	0.59
1:A:77:THR:HB	1:A:80:GLN:H	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:SER:OG	1:A:215:GLU:HG3	2.03	0.59
1:B:55:GLY:O	1:B:58:VAL:HG13	2.03	0.59
1:A:52:TRP:CZ3	1:A:160:SER:HB2	2.38	0.58
1:B:53:LEU:HD12	1:B:71:ASP:O	2.03	0.58
1:B:332:ASP:O	1:B:336:LEU:HD22	2.03	0.58
1:B:360:ARG:O	1:B:364:GLN:HG2	2.03	0.58
1:A:138:ASN:N	1:A:138:ASN:HD22	2.02	0.58
1:A:381:LEU:HD23	1:A:415:PHE:CE2	2.39	0.58
1:B:96:ARG:HG3	1:B:97:LEU:N	2.18	0.58
1:B:151:ASP:N	2:B:502:ATP:O2'	2.32	0.58
1:B:164:PHE:N	1:B:164:PHE:HD1	2.02	0.58
1:B:20:VAL:HG23	1:B:167:ARG:HD3	1.86	0.58
1:A:87:ASN:HD22	1:A:87:ASN:N	2.01	0.58
1:B:414:GLU:O	1:B:418:SER:HB3	2.04	0.58
1:A:337:ALA:O	1:A:341:VAL:HG23	2.04	0.57
1:A:26:HIS:O	1:A:28:ILE:N	2.37	0.57
1:A:225:THR:O	1:A:227:LEU:N	2.37	0.57
1:B:21:ILE:HB	1:B:166:VAL:HG12	1.85	0.57
1:A:299:MET:SD	1:B:378:TRP:HB3	2.45	0.57
1:A:269:THR:HA	1:A:272:PHE:CE2	2.39	0.57
1:B:223:LEU:O	1:B:226:LEU:HB2	2.05	0.57
1:A:220:ILE:HB	1:A:221:PRO:HD3	1.87	0.57
1:A:385:LYS:O	1:A:386:PHE:C	2.47	0.56
1:A:203:ARG:HB2	1:A:242:LEU:HD23	1.85	0.56
1:B:37:ALA:HB2	1:B:107:ILE:HG12	1.86	0.56
1:B:153:THR:HG22	1:B:170:VAL:HG12	1.87	0.56
1:B:395:LEU:O	1:B:396:ARG:C	2.48	0.56
1:A:22:PRO:HD2	1:A:25:GLN:CG	2.35	0.56
1:A:358:LEU:O	1:A:362:ILE:HG13	2.05	0.56
1:B:402:ASN:C	1:B:402:ASN:HD22	2.14	0.56
1:B:378:TRP:HZ3	1:B:423:GLN:OE1	1.88	0.56
1:B:270:ARG:HD2	1:B:271:TYR:HE2	1.68	0.56
1:B:152:PHE:N	1:B:152:PHE:CD2	2.73	0.56
1:B:27:ALA:O	1:B:28:ILE:C	2.49	0.56
1:B:34:SER:CB	1:B:38:LEU:HD21	2.36	0.56
1:B:302:ASN:ND2	1:B:303:PRO:HD2	2.20	0.56
1:B:402:ASN:C	1:B:402:ASN:ND2	2.63	0.55
1:A:196:VAL:HG23	1:A:230:ILE:HD13	1.89	0.55
1:A:212:ILE:HD13	1:A:258:TYR:OH	2.05	0.55
1:B:291:THR:O	1:B:295:ILE:HG12	2.06	0.55
1:B:152:PHE:CD1	1:B:181:ILE:HG22	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:LEU:O	1:B:230:ILE:HB	2.07	0.55
1:A:93:ARG:O	1:A:94:ARG:HB2	2.07	0.55
1:A:96:ARG:HH12	1:A:113:ARG:CD	2.02	0.55
1:A:402:ASN:HD22	1:A:402:ASN:C	2.12	0.55
1:A:170:VAL:HG12	1:A:171:GLY:N	2.20	0.54
1:A:238:GLU:O	1:A:242:LEU:HG	2.07	0.54
1:A:402:ASN:ND2	1:A:404:GLU:H	2.05	0.54
1:B:387:ARG:NH1	1:B:416:GLN:HE22	2.04	0.54
1:A:55:GLY:O	1:A:56:GLY:C	2.50	0.54
1:A:84:LEU:O	1:A:85:PHE:CD1	2.60	0.54
1:A:170:VAL:HG21	1:A:184:ILE:HD11	1.90	0.54
1:B:38:LEU:N	1:B:38:LEU:HD23	2.22	0.54
1:A:184:ILE:CG2	1:A:190:ARG:HH12	2.06	0.54
1:A:101:MET:HG2	1:A:102:PHE:N	2.17	0.54
1:A:244:GLN:HE22	1:A:303:PRO:CA	2.15	0.54
1:A:280:MET:HE2	1:A:314:PRO:CG	2.38	0.54
1:B:93:ARG:HG2	1:B:94:ARG:N	2.21	0.54
1:A:28:ILE:O	1:A:28:ILE:HG23	2.07	0.54
1:B:295:ILE:HD11	1:B:301:VAL:HG11	1.90	0.54
1:B:198:MET:HG2	1:B:220:ILE:HG12	1.91	0.53
1:B:370:SER:HA	1:B:408:LEU:HD11	1.91	0.53
1:A:23:ARG:HB2	1:A:164:PHE:HB3	1.90	0.53
1:A:181:ILE:CD1	1:A:210:MET:HG3	2.38	0.53
1:A:183:LEU:HB2	1:A:216:THR:CG2	2.38	0.53
1:A:198:MET:HG2	1:A:220:ILE:HG12	1.90	0.53
1:B:356:THR:O	1:B:360:ARG:HG2	2.08	0.53
1:A:86:ARG:HG3	1:A:87:ASN:ND2	2.18	0.53
1:A:26:HIS:CD2	1:A:26:HIS:C	2.86	0.53
1:B:170:VAL:CG1	1:B:171:GLY:H	2.20	0.53
1:B:368:ARG:HA	1:B:371:ARG:HD2	1.91	0.53
1:B:179:GLY:C	1:B:210:MET:HG2	2.34	0.53
1:B:405:LEU:O	1:B:406:GLN:C	2.51	0.53
1:A:109:VAL:HG12	1:A:109:VAL:O	2.08	0.53
1:B:244:GLN:OE1	1:B:306:LEU:HD22	2.08	0.53
1:A:22:PRO:HD2	1:A:25:GLN:CD	2.33	0.53
1:B:281:GLU:O	1:B:285:GLU:HG3	2.09	0.53
1:A:367:LEU:O	1:A:370:SER:HB3	2.09	0.53
1:B:311:PHE:C	1:B:314:PRO:HD2	2.34	0.53
1:B:405:LEU:C	1:B:407:ARG:N	2.64	0.53
1:B:418:SER:O	1:B:423:GLN:NE2	2.31	0.53
1:A:203:ARG:HB2	1:A:242:LEU:CD2	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:MET:HE2	1:A:314:PRO:HG3	1.90	0.53
1:A:18:VAL:HG22	1:A:18:VAL:O	2.08	0.52
1:A:40:VAL:HG11	1:A:107:ILE:HG21	1.90	0.52
1:A:184:ILE:HG23	1:A:185:GLY:N	2.24	0.52
1:B:59:ARG:NH1	2:B:502:ATP:O1G	2.41	0.52
1:B:73:THR:HB	1:B:112:PHE:CE2	2.44	0.52
1:B:150:ARG:HH21	1:B:150:ARG:HG3	1.74	0.52
1:B:323:ALA:O	1:B:327:GLY:HA2	2.10	0.52
1:A:223:LEU:O	1:A:226:LEU:HB2	2.09	0.52
1:B:154:ILE:HG12	1:B:155:ASN:N	2.23	0.52
1:B:325:GLU:O	1:B:326:SER:C	2.51	0.52
1:B:378:TRP:CZ3	1:B:423:GLN:OE1	2.62	0.52
1:B:23:ARG:HD2	1:B:164:PHE:HD2	1.75	0.52
1:A:94:ARG:HB3	1:A:97:LEU:HB3	1.91	0.52
1:A:94:ARG:NE	1:A:97:LEU:CD2	2.72	0.52
1:B:93:ARG:O	1:B:94:ARG:CB	2.56	0.52
1:B:303:PRO:HG2	1:B:355:LEU:HD11	1.92	0.52
1:A:234:HIS:ND1	1:A:234:HIS:N	2.55	0.52
1:A:396:ARG:O	1:A:400:GLU:HG3	2.09	0.52
1:B:153:THR:HB	1:B:172:GLY:H	1.75	0.51
1:A:396:ARG:NH2	1:A:405:LEU:HD11	2.24	0.51
1:A:150:ARG:HH21	1:A:150:ARG:HG3	1.76	0.51
1:B:41:MET:HB3	1:B:51:ALA:HB1	1.91	0.51
1:A:365:LEU:HD12	1:A:389:ALA:HB2	1.92	0.51
1:B:199:LEU:O	1:B:202:VAL:HG23	2.10	0.51
1:A:268:ILE:O	1:A:269:THR:C	2.53	0.51
1:A:278:SER:O	1:A:282:ARG:HD3	2.10	0.51
1:A:383:HIS:CG	1:A:384:PRO:HD2	2.45	0.51
1:A:208:LEU:CD1	1:A:210:MET:HE2	2.37	0.51
1:B:23:ARG:HA	1:B:26:HIS:CE1	2.45	0.51
1:A:297:ASN:ND2	1:B:378:TRP:NE1	2.57	0.51
1:B:113:ARG:HD3	1:B:146:ASP:OD1	2.11	0.51
1:B:261:PHE:CE2	1:B:265:PHE:HB2	2.45	0.51
1:A:417:VAL:HG12	1:B:416:GLN:NE2	2.26	0.50
1:B:291:THR:HG23	1:B:301:VAL:HG21	1.93	0.50
1:B:158:TYR:O	1:B:166:VAL:HG23	2.11	0.50
1:B:183:LEU:CD2	1:B:184:ILE:H	2.10	0.50
1:B:305:PHE:O	1:B:306:LEU:C	2.53	0.50
1:B:415:PHE:O	1:B:415:PHE:CD2	2.64	0.50
1:B:304:ALA:HB2	1:B:355:LEU:HD12	1.92	0.50
1:A:150:ARG:HD2	1:A:156:SER:OG	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:TYR:CE1	1:B:317:GLU:HG2	2.44	0.50
1:B:325:GLU:HG2	1:B:326:SER:N	2.26	0.50
1:A:27:ALA:C	1:A:29:SER:H	2.19	0.50
1:A:143:ILE:HD13	1:A:143:ILE:O	2.11	0.50
1:A:402:ASN:HD21	1:A:404:GLU:HB3	1.76	0.50
1:B:233:ALA:HA	1:B:350:ALA:HB3	1.94	0.50
1:A:358:LEU:HD23	1:A:385:LYS:CE	2.41	0.50
1:B:302:ASN:O	1:B:305:PHE:HB3	2.12	0.50
1:A:35:GLU:O	1:A:37:ALA:N	2.41	0.50
1:A:181:ILE:HD11	1:A:210:MET:HG3	1.92	0.50
1:A:413:GLY:O	1:B:417:VAL:HG11	2.11	0.50
1:B:381:LEU:HD22	1:B:415:PHE:HE2	1.77	0.50
1:A:410:LYS:O	1:A:414:GLU:HG3	2.11	0.49
1:B:26:HIS:C	1:B:26:HIS:CD2	2.90	0.49
1:A:113:ARG:HG3	1:A:149:ARG:HG3	1.94	0.49
1:A:218:GLU:O	1:A:221:PRO:HD2	2.11	0.49
1:B:280:MET:O	1:B:284:ILE:HG13	2.13	0.49
1:B:219:PRO:CB	1:B:223:LEU:HD12	2.43	0.49
1:B:244:GLN:NE2	1:B:303:PRO:CA	2.76	0.49
1:B:197:ARG:NH1	2:B:502:ATP:N1	2.60	0.49
1:B:213:SER:HB3	1:B:216:THR:OG1	2.11	0.49
1:A:61:LEU:HG	1:A:67:PRO:HG3	1.95	0.49
1:A:202:VAL:HG11	1:A:255:LEU:HD21	1.95	0.49
1:B:23:ARG:HB3	1:B:164:PHE:HB3	1.94	0.49
1:B:30:ARG:HE	1:B:33:ILE:CG1	2.26	0.49
1:A:28:ILE:HD12	1:A:30:ARG:HG2	1.95	0.48
1:B:73:THR:HA	1:B:110:ALA:O	2.13	0.48
1:A:202:VAL:HG11	1:A:255:LEU:CD2	2.43	0.48
1:B:261:PHE:C	1:B:261:PHE:CD2	2.91	0.48
1:A:385:LYS:O	1:A:387:ARG:N	2.47	0.48
1:B:283:ILE:CG2	1:B:284:ILE:N	2.76	0.48
1:B:184:ILE:HG22	1:B:185:GLY:N	2.28	0.48
1:A:174:LYS:HE2	1:A:178:ASP:OD2	2.12	0.48
1:B:26:HIS:O	1:B:26:HIS:CD2	2.67	0.48
1:A:85:PHE:O	1:A:88:CYS:HB2	2.14	0.48
1:A:20:VAL:HG23	1:A:167:ARG:HD3	1.96	0.48
1:A:251:THR:HG22	1:A:251:THR:O	2.14	0.48
1:A:414:GLU:CB	1:A:426:MET:HE1	2.44	0.48
1:A:187:PRO:HB3	1:A:191:TYR:HE2	1.79	0.48
1:A:184:ILE:HG22	1:A:185:GLY:H	1.79	0.47
1:B:50:GLU:HG2	1:B:52:TRP:HE1	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LYS:HA	1:B:67:PRO:HD2	1.61	0.47
1:B:248:GLY:C	1:B:288:LEU:HD13	2.39	0.47
1:B:289:LYS:O	1:B:292:ASP:HB2	2.14	0.47
1:A:72:VAL:CG2	1:A:109:VAL:HG22	2.43	0.47
1:A:94:ARG:CZ	1:A:97:LEU:HD21	2.43	0.47
1:A:280:MET:HG3	1:A:313:TYR:HB2	1.95	0.47
1:A:419:ALA:HB1	1:A:421:PRO:HD2	1.97	0.47
1:A:58:VAL:O	1:A:62:LEU:HD12	2.15	0.47
1:A:152:PHE:N	1:A:152:PHE:CD2	2.83	0.47
1:A:179:GLY:C	1:A:210:MET:HB2	2.39	0.47
1:B:290:ASN:C	1:B:292:ASP:N	2.68	0.47
1:A:246:GLY:HA2	1:A:292:ASP:HA	1.96	0.47
1:A:305:PHE:O	1:A:306:LEU:C	2.57	0.47
1:A:391:ASP:O	1:A:394:ALA:HB3	2.15	0.47
1:A:312:TRP:CE2	1:A:316:LEU:HD22	2.50	0.47
1:A:21:ILE:HA	1:A:22:PRO:HD3	1.48	0.47
1:A:94:ARG:HG3	1:A:97:LEU:HD23	1.97	0.47
1:A:195:PRO:O	1:A:198:MET:HB2	2.15	0.47
1:B:149:ARG:O	2:B:502:ATP:C2	2.68	0.47
1:B:280:MET:HE2	1:B:314:PRO:CG	2.45	0.47
1:B:52:TRP:HA	1:B:159:TYR:O	2.15	0.46
1:A:371:ARG:O	1:A:373:GLN:N	2.49	0.46
1:B:158:TYR:O	1:B:166:VAL:HA	2.15	0.46
1:A:243:LEU:HA	1:A:248:GLY:HA2	1.98	0.46
1:B:224:ALA:O	1:B:225:THR:C	2.58	0.46
1:A:97:LEU:O	1:A:97:LEU:HG	2.16	0.46
1:B:224:ALA:O	1:B:226:LEU:N	2.48	0.46
1:A:278:SER:HB2	1:A:279:PRO:HD2	1.98	0.46
1:B:182:ARG:HG2	1:B:183:LEU:O	2.15	0.46
1:B:359:THR:O	1:B:360:ARG:C	2.56	0.46
1:B:280:MET:HE2	1:B:314:PRO:HG3	1.98	0.46
1:A:171:GLY:O	1:A:172:GLY:C	2.59	0.46
1:B:72:VAL:HG23	1:B:109:VAL:HG22	1.98	0.46
1:B:268:ILE:HD11	1:B:310:MET:HB3	1.97	0.46
1:A:387:ARG:HA	1:A:387:ARG:HD3	1.59	0.46
1:B:387:ARG:HH12	1:B:416:GLN:HE22	1.63	0.46
1:A:195:PRO:HB2	1:A:227:LEU:HD23	1.97	0.45
1:A:307:PHE:O	1:A:308:ALA:C	2.55	0.45
1:B:30:ARG:NE	1:B:33:ILE:HD11	2.31	0.45
1:B:56:GLY:O	1:B:59:ARG:HB3	2.17	0.45
1:B:212:ILE:HG22	1:B:213:SER:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:LEU:C	1:B:263:PRO:HD2	2.41	0.45
1:A:52:TRP:CE3	1:A:160:SER:HB2	2.51	0.45
1:A:158:TYR:O	1:A:166:VAL:HA	2.17	0.45
1:B:260:LEU:O	1:B:263:PRO:HD2	2.17	0.45
1:B:278:SER:HB2	1:B:279:PRO:CD	2.46	0.45
1:A:400:GLU:O	1:A:401:ARG:C	2.57	0.45
1:A:280:MET:O	1:A:284:ILE:HG13	2.17	0.45
1:A:359:THR:O	1:A:360:ARG:C	2.58	0.45
1:B:345:ALA:C	1:B:347:ARG:H	2.23	0.45
1:B:236:PHE:HA	1:B:349:LEU:HD12	1.99	0.45
1:A:223:LEU:HD23	1:A:226:LEU:CD1	2.47	0.44
1:A:381:LEU:HD23	1:A:415:PHE:CD2	2.51	0.44
1:B:247:TYR:O	1:B:249:TYR:N	2.50	0.44
1:A:75:ASN:C	1:A:75:ASN:ND2	2.75	0.44
1:A:143:ILE:HD13	1:A:143:ILE:C	2.43	0.44
1:A:152:PHE:CD1	1:A:181:ILE:HG23	2.53	0.44
1:A:219:PRO:HB3	1:A:223:LEU:CD1	2.48	0.44
1:A:312:TRP:NE1	1:A:316:LEU:HD22	2.32	0.44
1:A:84:LEU:HD23	1:A:85:PHE:CE1	2.52	0.44
1:A:146:ASP:C	1:A:148:GLN:H	2.25	0.44
1:A:402:ASN:C	1:A:402:ASN:ND2	2.75	0.44
1:B:21:ILE:HG13	1:B:166:VAL:HG13	1.99	0.44
1:B:103:GLY:C	1:B:105:GLU:H	2.25	0.44
1:B:266:PRO:O	1:B:268:ILE:N	2.50	0.44
1:A:408:LEU:N	1:A:408:LEU:HD23	2.31	0.44
1:A:427:LEU:HD23	1:A:427:LEU:HA	1.77	0.44
1:B:186:ASN:O	1:B:187:PRO:C	2.60	0.44
1:B:232:PRO:CB	1:B:350:ALA:HB2	2.44	0.44
1:A:402:ASN:O	1:A:406:GLN:HB2	2.17	0.44
1:B:244:GLN:HE22	1:B:303:PRO:CA	2.27	0.44
1:A:83:LYS:HG2	1:A:86:ARG:HH21	1.83	0.44
1:B:36:ASN:O	1:B:102:PHE:CZ	2.71	0.44
1:B:233:ALA:O	1:B:234:HIS:C	2.58	0.44
1:A:224:ALA:O	1:A:225:THR:C	2.57	0.44
1:A:289:LYS:O	1:A:292:ASP:HB2	2.18	0.44
1:A:85:PHE:CD1	1:A:85:PHE:N	2.86	0.43
1:A:152:PHE:CZ	1:A:201:ALA:HA	2.53	0.43
1:B:41:MET:HB3	1:B:51:ALA:CB	2.48	0.43
1:B:153:THR:HG22	1:B:170:VAL:HG11	1.99	0.43
1:B:280:MET:O	1:B:280:MET:HG2	2.16	0.43
1:B:364:GLN:HE21	1:B:364:GLN:HA	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:368:ARG:NH1	1:A:380:LEU:HD11	2.33	0.43
1:B:235:LEU:O	1:B:236:PHE:C	2.61	0.43
1:B:278:SER:O	1:B:279:PRO:C	2.57	0.43
1:A:20:VAL:O	1:A:20:VAL:HG13	2.17	0.43
1:A:235:LEU:O	1:A:239:SER:HB3	2.18	0.43
1:B:59:ARG:NE	1:B:208:LEU:HD21	2.33	0.43
1:A:106:ILE:O	1:A:107:ILE:HD13	2.18	0.43
1:A:361:ASP:O	1:A:365:LEU:HG	2.19	0.43
1:A:61:LEU:HD23	1:A:61:LEU:HA	1.61	0.43
1:A:79:GLU:HG3	1:A:82:ARG:HH21	1.83	0.43
1:A:382:GLU:O	1:A:383:HIS:C	2.61	0.43
1:A:418:SER:O	1:A:423:GLN:NE2	2.45	0.43
1:B:144:GLU:HA	1:B:169:TYR:CE2	2.53	0.43
1:B:268:ILE:HD13	1:B:284:ILE:HD11	2.01	0.43
1:A:37:ALA:O	1:A:40:VAL:CG1	2.65	0.43
1:A:225:THR:C	1:A:227:LEU:H	2.25	0.43
1:B:185:GLY:H	1:B:190:ARG:NH1	2.17	0.43
1:B:288:LEU:O	1:B:289:LYS:C	2.61	0.43
1:B:339:ASN:O	1:B:343:ASP:HB2	2.18	0.43
1:A:28:ILE:CG2	1:A:30:ARG:CG	2.72	0.43
1:A:94:ARG:HB3	1:A:97:LEU:CD2	2.40	0.43
1:A:414:GLU:HB3	1:A:426:MET:HE1	2.00	0.43
1:A:151:ASP:OD1	1:A:151:ASP:N	2.51	0.43
1:A:368:ARG:CZ	1:A:380:LEU:HD11	2.49	0.43
1:A:416:GLN:NE2	1:B:417:VAL:HG12	2.34	0.43
1:A:415:PHE:CD2	1:A:415:PHE:O	2.72	0.43
1:B:264:LEU:O	1:B:266:PRO:HD2	2.18	0.43
1:B:268:ILE:CD1	1:B:310:MET:HB3	2.49	0.42
1:A:86:ARG:CG	1:A:87:ASN:HD22	2.25	0.42
1:A:230:ILE:HG23	1:A:231:PRO:HD2	2.02	0.42
1:A:380:LEU:C	1:A:382:GLU:N	2.75	0.42
1:A:421:PRO:HD3	1:B:290:ASN:ND2	2.34	0.42
1:B:251:THR:O	1:B:255:LEU:HG	2.19	0.42
1:B:315:LEU:HA	1:B:341:VAL:HG21	2.00	0.42
1:B:406:GLN:O	1:B:409:VAL:HB	2.18	0.42
1:A:271:TYR:HB2	1:A:280:MET:HE1	2.00	0.42
1:B:238:GLU:O	1:B:242:LEU:HD12	2.19	0.42
1:B:295:ILE:HD11	1:B:301:VAL:CG1	2.48	0.42
1:A:182:ARG:HH11	1:A:182:ARG:HG3	1.85	0.42
1:B:95:PHE:C	1:B:96:ARG:HG2	2.43	0.42
1:B:219:PRO:O	1:B:223:LEU:HD12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ILE:HG13	1:A:155:ASN:N	2.34	0.42
1:A:383:HIS:ND1	1:A:384:PRO:HD2	2.34	0.42
1:A:395:LEU:O	1:A:396:ARG:C	2.62	0.42
1:B:208:LEU:HD12	1:B:208:LEU:H	1.84	0.42
1:B:364:GLN:HA	1:B:364:GLN:NE2	2.34	0.42
1:B:402:ASN:ND2	1:B:404:GLU:H	2.18	0.42
1:A:225:THR:C	1:A:227:LEU:N	2.77	0.42
1:A:264:LEU:C	1:A:266:PRO:HD3	2.44	0.42
1:A:280:MET:HA	1:A:313:TYR:CD2	2.53	0.42
1:B:138:ASN:OD1	1:B:138:ASN:N	2.52	0.42
1:B:170:VAL:CG1	1:B:171:GLY:N	2.75	0.42
1:B:304:ALA:HB2	1:B:355:LEU:CD1	2.50	0.42
1:B:386:PHE:O	1:B:387:ARG:C	2.63	0.42
1:A:269:THR:HA	1:A:272:PHE:CD2	2.54	0.42
1:B:103:GLY:O	1:B:105:GLU:N	2.53	0.42
1:B:154:ILE:C	1:B:156:SER:H	2.28	0.42
1:B:296:HIS:C	1:B:298:ASP:H	2.26	0.42
1:A:150:ARG:HG3	1:A:150:ARG:NH2	2.34	0.42
1:A:196:VAL:C	1:A:198:MET:N	2.76	0.42
1:B:316:LEU:O	1:B:319:ALA:HB3	2.20	0.42
1:A:54:VAL:HG22	1:A:73:THR:HG22	2.01	0.42
1:A:375:LYS:HD2	1:A:375:LYS:H	1.85	0.42
1:B:94:ARG:HD2	1:B:94:ARG:HA	2.00	0.42
1:B:275:ASN:ND2	1:B:277:ASP:H	2.18	0.42
1:B:283:ILE:HG23	1:B:284:ILE:N	2.35	0.42
1:B:354:ARG:NH1	1:B:355:LEU:HD21	2.35	0.42
1:A:402:ASN:ND2	1:A:404:GLU:N	2.67	0.42
1:B:402:ASN:HD21	1:B:404:GLU:HB3	1.85	0.42
1:B:405:LEU:HD23	1:B:405:LEU:HA	1.75	0.42
1:A:280:MET:CE	1:A:314:PRO:HG3	2.50	0.41
1:B:89:ARG:NH1	1:B:101:MET:HE2	2.35	0.41
1:B:96:ARG:O	1:B:97:LEU:C	2.62	0.41
1:B:103:GLY:C	1:B:105:GLU:N	2.78	0.41
1:B:186:ASN:ND2	1:B:188:GLU:HB3	2.35	0.41
1:A:184:ILE:HG22	1:A:185:GLY:N	2.35	0.41
1:B:152:PHE:H	1:B:152:PHE:HD2	1.67	0.41
1:B:293:THR:O	1:B:294:ARG:C	2.62	0.41
1:B:391:ASP:O	1:B:394:ALA:HB3	2.20	0.41
1:A:77:THR:C	1:A:79:GLU:N	2.77	0.41
1:B:45:ASN:HD22	1:B:45:ASN:HA	1.66	0.41
1:A:35:GLU:C	1:A:37:ALA:H	2.26	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ALA:CB	1:A:107:ILE:HG13	2.50	0.41
1:A:152:PHE:CZ	1:A:201:ALA:HB2	2.56	0.41
1:A:315:LEU:HD22	1:A:363:TRP:CE3	2.56	0.41
1:A:364:GLN:HE21	1:A:364:GLN:HA	1.85	0.41
1:B:61:LEU:HD23	1:B:67:PRO:HD3	2.03	0.41
1:A:191:TYR:CE1	1:A:198:MET:CG	2.99	0.41
1:A:385:LYS:O	1:A:388:ALA:N	2.54	0.41
1:B:37:ALA:HB2	1:B:107:ILE:CG1	2.49	0.41
1:A:30:ARG:C	1:A:32:ASP:N	2.79	0.41
1:A:247:TYR:O	1:A:248:GLY:C	2.64	0.41
1:A:286:GLN:NE2	1:A:286:GLN:HA	2.35	0.41
1:A:351:ILE:HA	1:A:352:PRO:HD3	1.87	0.41
1:A:372:ARG:NH1	1:A:408:LEU:CD2	2.84	0.41
1:B:202:VAL:O	1:B:205:ALA:HB3	2.20	0.41
1:A:372:ARG:HD2	1:A:372:ARG:HA	1.80	0.41
1:B:266:PRO:C	1:B:268:ILE:N	2.79	0.41
1:A:194:ASP:HB3	1:A:197:ARG:HG3	2.03	0.40
1:A:378:TRP:CE2	1:B:297:ASN:ND2	2.89	0.40
1:B:67:PRO:HG3	1:B:70:PHE:CZ	2.55	0.40
1:A:316:LEU:HD12	1:A:316:LEU:HA	1.76	0.40
1:A:365:LEU:HD12	1:A:389:ALA:CB	2.51	0.40
1:A:413:GLY:O	1:A:416:GLN:HG2	2.22	0.40
1:B:96:ARG:CG	1:B:97:LEU:N	2.82	0.40
1:B:172:GLY:O	1:B:173:MET:C	2.64	0.40
1:B:184:ILE:HG22	1:B:190:ARG:HH12	1.84	0.40
1:B:224:ALA:C	1:B:226:LEU:N	2.79	0.40
1:B:235:LEU:HD13	1:B:264:LEU:HD22	2.02	0.40
1:A:196:VAL:CG2	1:A:230:ILE:HD13	2.49	0.40
1:A:90:LEU:CD2	1:A:90:LEU:H	2.35	0.40
1:A:146:ASP:C	1:A:148:GLN:N	2.80	0.40
1:A:203:ARG:HD3	1:A:242:LEU:HD23	2.02	0.40
1:A:302:ASN:HD21	1:A:385:LYS:HE3	1.87	0.40
1:B:247:TYR:O	1:B:248:GLY:C	2.64	0.40
1:B:218:GLU:N	1:B:219:PRO:CD	2.85	0.40
1:B:337:ALA:O	1:B:341:VAL:HG23	2.22	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	384/415 (92%)	334 (87%)	41 (11%)	9 (2%)	5	25
1	B	384/415 (92%)	333 (87%)	40 (10%)	11 (3%)	3	21
All	All	768/830 (92%)	667 (87%)	81 (10%)	20 (3%)	4	23

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ILE
1	A	248	GLY
1	B	93	ARG
1	B	94	ARG
1	B	97	LEU
1	B	104	PRO
1	B	248	GLY
1	A	172	GLY
1	A	226	LEU
1	B	151	ASP
1	A	29	SER
1	B	172	GLY
1	B	327	GLY
1	B	267	THR
1	A	112	PHE
1	A	421	PRO
1	B	56	GLY
1	B	92	GLY
1	A	56	GLY
1	A	327	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	334/358 (93%)	278 (83%)	56 (17%)	2	11
1	B	334/358 (93%)	272 (81%)	62 (19%)	1	7
All	All	668/716 (93%)	550 (82%)	118 (18%)	2	8

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	HIS
1	A	28	ILE
1	A	29	SER
1	A	32	ASP
1	A	40	VAL
1	A	41	MET
1	A	53	LEU
1	A	54	VAL
1	A	71	ASP
1	A	72	VAL
1	A	73	THR
1	A	75	ASN
1	A	82	ARG
1	A	84	LEU
1	A	87	ASN
1	A	90	LEU
1	A	93	ARG
1	A	100	VAL
1	A	101	MET
1	A	106	ILE
1	A	138	ASN
1	A	140	PHE
1	A	142	SER
1	A	143	ILE
1	A	145	GLU
1	A	149	ARG
1	A	151	ASP
1	A	154	ILE
1	A	156	SER
1	A	161	VAL
1	A	166	VAL

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Mol	Chain	Res	Type
1	A	176	LEU
1	A	184	ILE
1	A	208	LEU
1	A	210	MET
1	A	212	ILE
1	A	215	GLU
1	A	226	LEU
1	A	234	HIS
1	A	256	CYS
1	A	262	GLN
1	A	277	ASP
1	A	280	MET
1	A	283	ILE
1	A	286	GLN
1	A	291	THR
1	A	329	THR
1	A	349	LEU
1	A	351	ILE
1	A	355	LEU
1	A	364	GLN
1	A	375	LYS
1	A	381	LEU
1	A	387	ARG
1	A	402	ASN
1	A	405	LEU
1	B	20	VAL
1	B	21	ILE
1	B	24	GLU
1	B	26	HIS
1	B	28	ILE
1	B	30	ARG
1	B	33	ILE
1	B	35	GLU
1	B	38	LEU
1	B	50	GLU
1	B	54	VAL
1	B	58	VAL
1	B	71	ASP
1	B	73	THR
1	B	77	THR
1	B	100	VAL
1	B	104	PRO

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Mol	Chain	Res	Type
1	B	107	ILE
1	B	113	ARG
1	B	138	ASN
1	B	139	ILE
1	B	152	PHE
1	B	154	ILE
1	B	156	SER
1	B	163	ASP
1	B	164	PHE
1	B	165	THR
1	B	184	ILE
1	B	189	THR
1	B	202	VAL
1	B	208	LEU
1	B	210	MET
1	B	215	GLU
1	B	225	THR
1	B	226	LEU
1	B	254	LEU
1	B	268	ILE
1	B	269	THR
1	B	270	ARG
1	B	275	ASN
1	B	277	ASP
1	B	283	ILE
1	B	284	ILE
1	B	288	LEU
1	B	298	ASP
1	B	300	ARG
1	B	316	LEU
1	B	324	GLN
1	B	329	THR
1	B	336	LEU
1	B	342	LEU
1	B	343	ASP
1	B	346	CYS
1	B	371	ARG
1	B	372	ARG
1	B	380	LEU
1	B	385	LYS
1	B	393	LEU
1	B	402	ASN

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Mol	Chain	Res	Type
1	B	405	LEU
1	B	406	GLN
1	B	418	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	26	HIS
1	A	45	ASN
1	A	75	ASN
1	A	80	GLN
1	A	87	ASN
1	A	155	ASN
1	A	244	GLN
1	A	275	ASN
1	A	290	ASN
1	A	302	ASN
1	A	364	GLN
1	A	402	ASN
1	A	406	GLN
1	A	416	GLN
1	B	17	GLN
1	B	26	HIS
1	B	45	ASN
1	B	87	ASN
1	B	155	ASN
1	B	186	ASN
1	B	244	GLN
1	B	275	ASN
1	B	290	ASN
1	B	302	ASN
1	B	364	GLN
1	B	373	GLN
1	B	402	ASN
1	B	416	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ATP	B	502	3	32,33,33	1.84	6 (18%)	48,52,52	1.84	8 (16%)
2	ATP	A	501	3	32,33,33	1.84	6 (18%)	48,52,52	1.84	8 (16%)

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	3	-	1/22/38/38	0/3/3/3
2	ATP	A	501	3	-	1/22/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	ATP	C5-C4	5.32	1.48	1.39
2	B	502	ATP	C5-C4	5.28	1.48	1.39
2	B	502	ATP	PB-O3B	4.16	1.64	1.59
2	A	501	ATP	PB-O3B	4.12	1.63	1.59
2	A	501	ATP	PB-O3A	3.99	1.63	1.59
2	B	502	ATP	PB-O3A	3.97	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	ATP	PA-O3A	3.59	1.63	1.59
2	A	501	ATP	PA-O3A	3.58	1.63	1.59
2	B	502	ATP	C5-C6	3.36	1.50	1.41
2	A	501	ATP	C5-C6	3.33	1.50	1.41
2	A	501	ATP	C8-N7	2.50	1.36	1.31
2	B	502	ATP	C8-N7	2.48	1.36	1.31

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ATP	C5-C4-N3	-6.63	117.58	126.72
2	B	502	ATP	C5-C4-N3	-6.60	117.62	126.72
2	B	502	ATP	N3-C4-N9	4.97	135.62	127.17
2	A	501	ATP	N3-C4-N9	4.95	135.58	127.17
2	B	502	ATP	C2-N3-C4	4.10	121.85	111.83
2	A	501	ATP	C2-N3-C4	4.09	121.81	111.83
2	A	501	ATP	C4-C5-N7	-3.74	106.31	110.58
2	B	502	ATP	C4-C5-N7	-3.69	106.36	110.58
2	A	501	ATP	N3-C2-N1	-3.24	123.68	128.58
2	B	502	ATP	N3-C2-N1	-3.24	123.68	128.58
2	A	501	ATP	C3'-C2'-C1'	2.56	106.30	101.46
2	B	502	ATP	C3'-C2'-C1'	2.55	106.29	101.46
2	B	502	ATP	C5-N7-C8	2.42	107.25	103.45
2	A	501	ATP	C5-N7-C8	2.41	107.24	103.45
2	A	501	ATP	C6-C5-N7	2.08	136.11	132.09
2	B	502	ATP	C6-C5-N7	2.07	136.07	132.09

There are no chirality outliers.

All (2) torsion outliers are listed below:

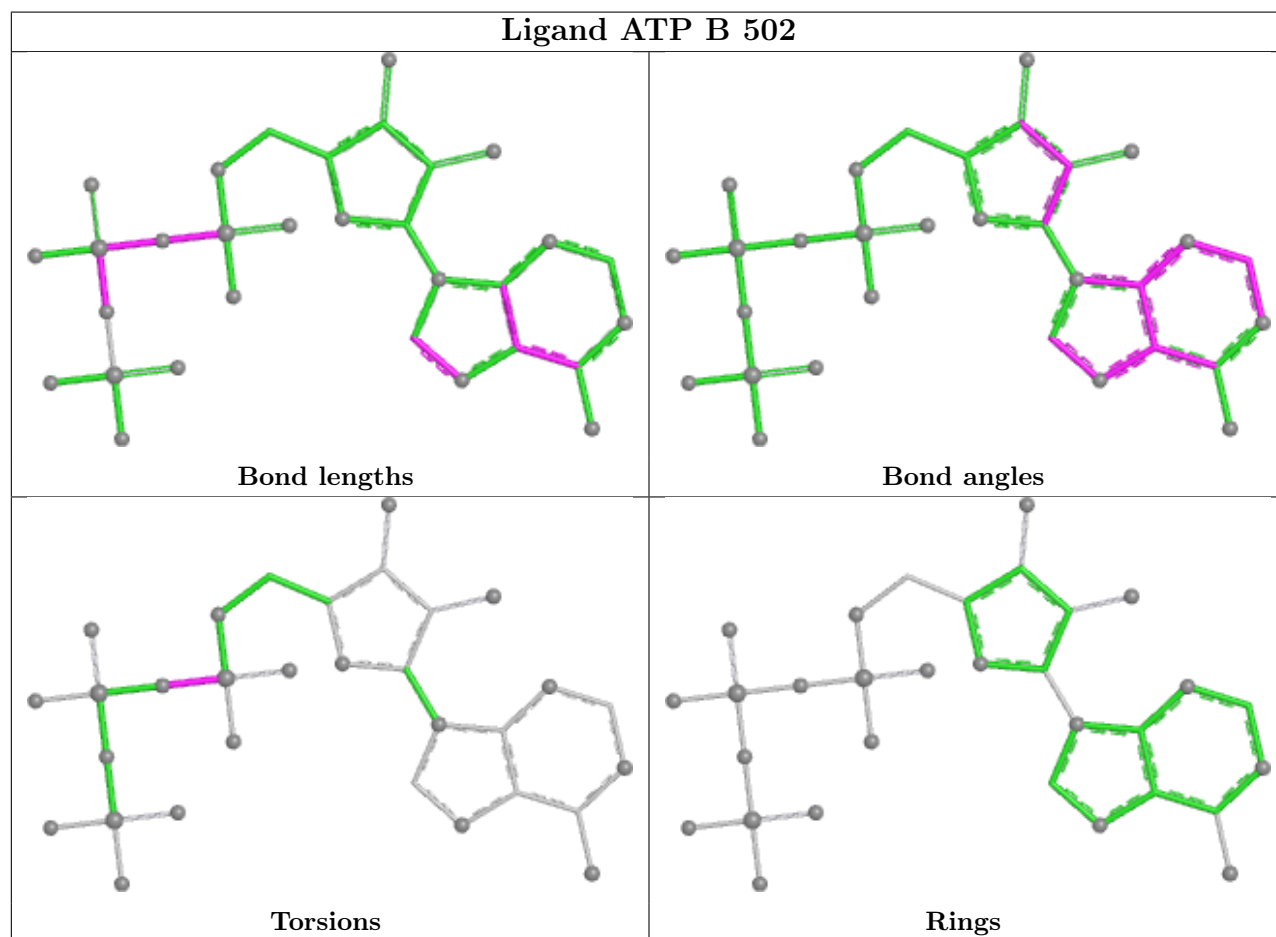
Mol	Chain	Res	Type	Atoms
2	A	501	ATP	PB-O3A-PA-O5'
2	B	502	ATP	PB-O3A-PA-O5'

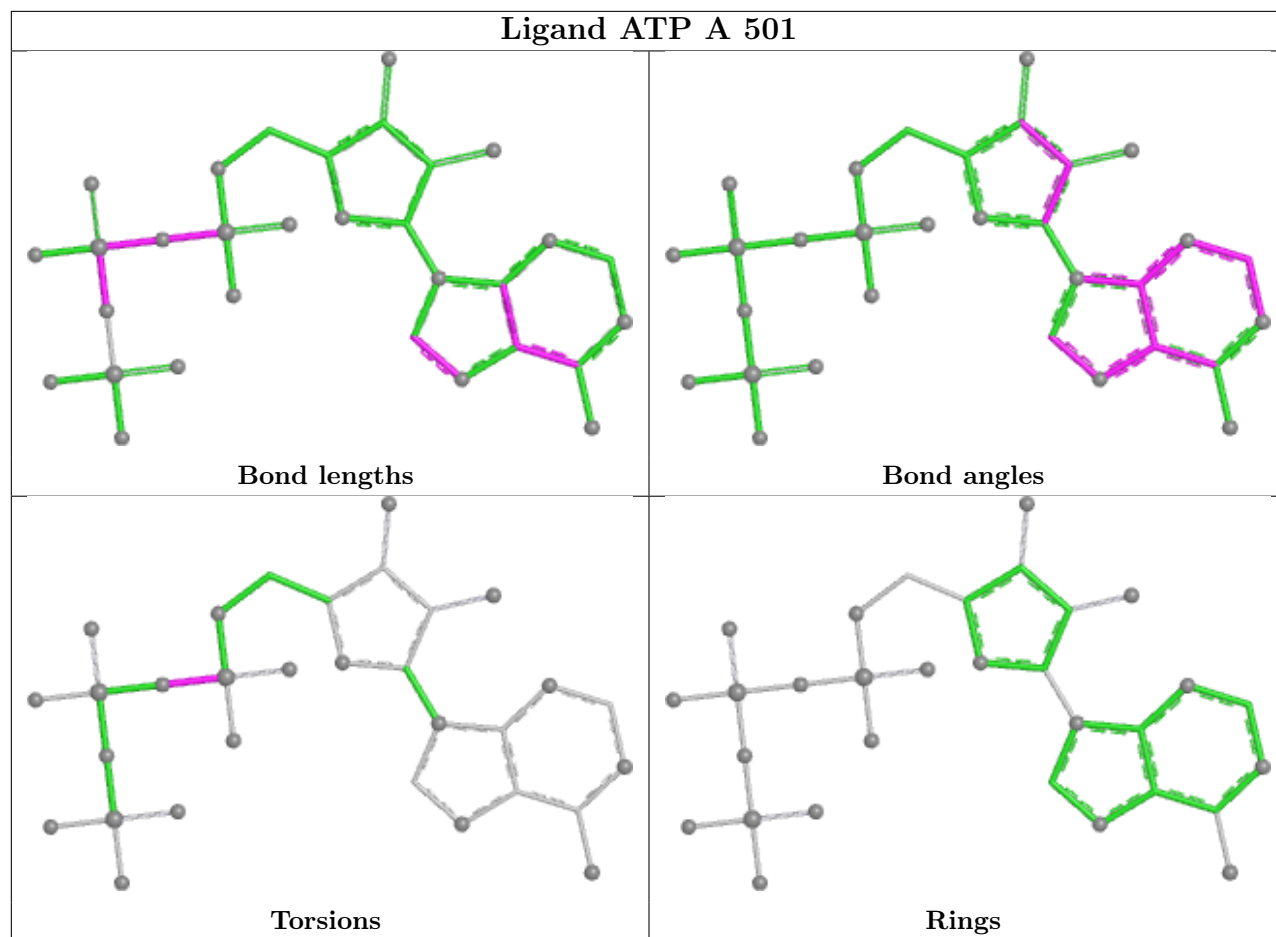
There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	ATP	7	0
2	A	501	ATP	4	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 [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	388/415 (93%)	-0.43	0 100 100	75, 122, 169, 219	0
1	B	388/415 (93%)	-0.47	0 100 100	90, 123, 170, 205	0
All	All	776/830 (93%)	-0.45	0 100 100	75, 123, 170, 219	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

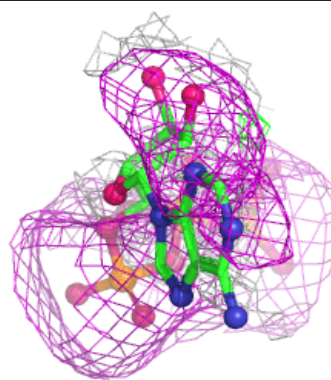
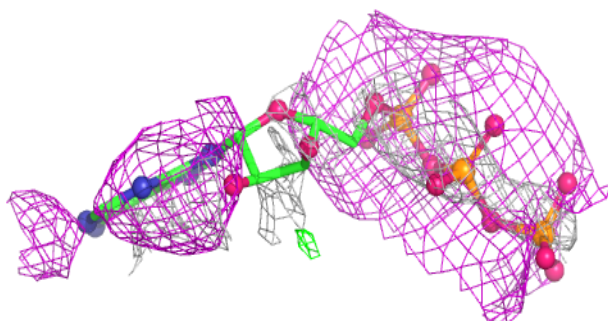
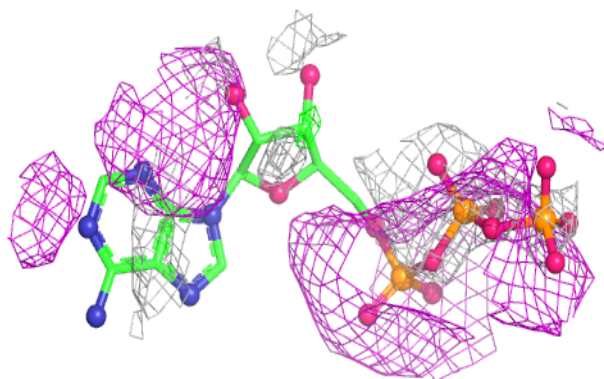
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ATP	B	502	31/31	0.53	0.17	240,245,246,248	0
2	ATP	A	501	31/31	0.56	0.17	234,236,242,242	0
3	MG	A	600	1/1	0.88	0.07	117,117,117,117	0
3	MG	B	601	1/1	0.91	0.23	93,93,93,93	0

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.

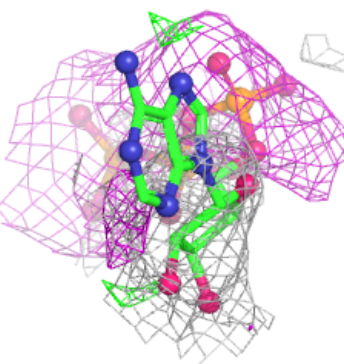
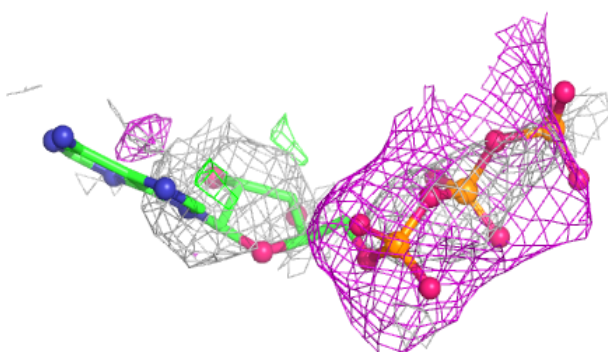
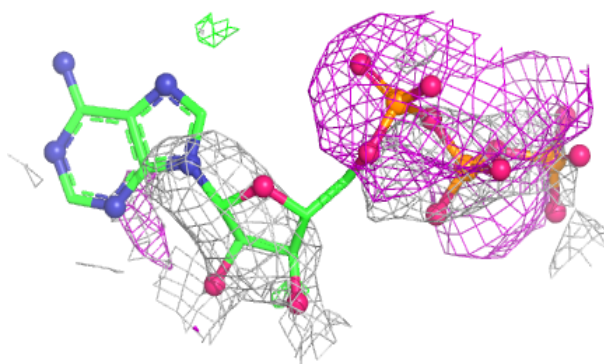
Electron density around ATP B 502:

$2mF_o-DF_c$ (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.