



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:04 PM UTC

PDB ID : 4AT8 / pdb_00004at8
Title : Crystal structure of the NF90-NF45 dimerisation domain complex with ATP
Authors : Wolkowicz, U.M.; Cook, A.G.
Deposited on : 2012-05-05
Resolution : 2.69 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

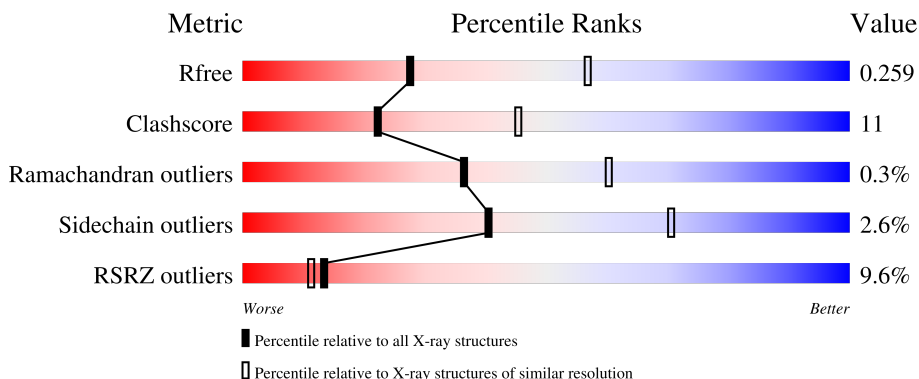
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	364	8% 70% 20% • 9%
1	C	364	9% 70% 20% • 9%
2	B	383	10% 65% 16% • 17%
2	D	383	6% 69% 13% • 17%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9798 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 INTERLEUKIN ENHANCER-BINDING FACTOR 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2538	1629	423	476	10			
1	C	332	Total	C	N	O	S	0	0	0
			2495	1596	418	471	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9CXY6
A	0	SER	-	expression tag	UNP Q9CXY6
C	27	GLY	-	expression tag	UNP Q9CXY6
C	28	SER	-	expression tag	UNP Q9CXY6

- Molecule 2 is a protein called INTERLEUKIN ENHANCER-BINDING FACTOR 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	316	Total	C	N	O	S	0	0	0
			2332	1469	409	441	13			
2	D	316	Total	C	N	O	S	0	0	0
			2369	1498	407	450	14			

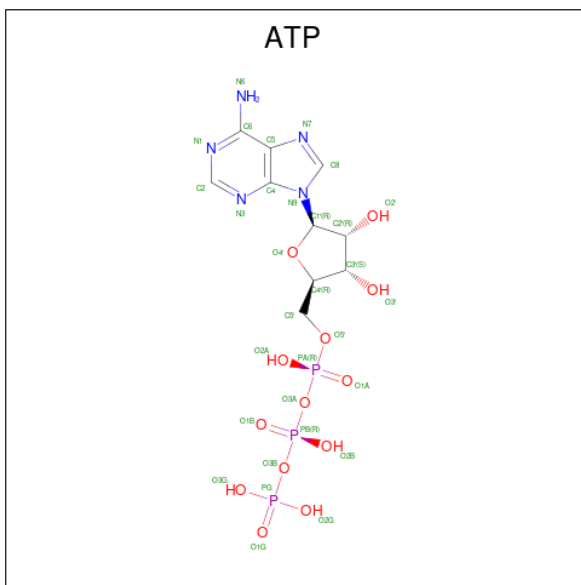
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP Q9Z1X4
B	0	HIS	-	expression tag	UNP Q9Z1X4
D	-1	GLY	-	expression tag	UNP Q9Z1X4
D	0	HIS	-	expression tag	UNP Q9Z1X4

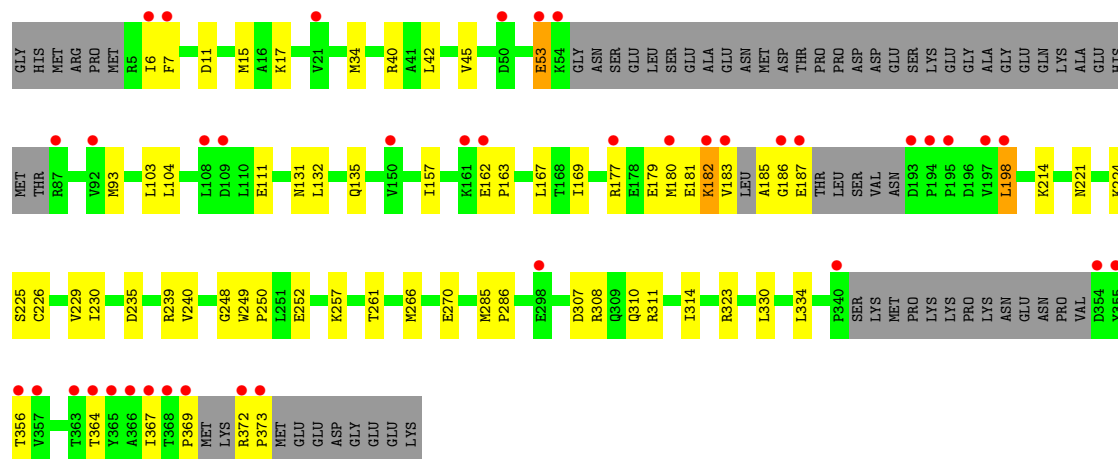
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $\text{C}_{10}\text{H}_{16}\text{N}_5\text{O}_{13}\text{P}_3$).



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



• Molecule 2: INTERLEUKIN ENHANCER-BINDING FACTOR 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.87Å 84.13Å 159.60Å 90.00° 93.34° 90.00°	Depositor
Resolution (Å)	57.84 – 2.69 57.84 – 2.69	Depositor EDS
% Data completeness (in resolution range)	96.9 (57.84-2.69) 96.7 (57.84-2.69)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.240 , 0.266 0.234 , 0.259	Depositor DCC
R_{free} test set	2574 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.2	Xtriage
Anisotropy	0.564	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.1	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	0.90	EDS
Total number of atoms	9798	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.20% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.37	0/2588	0.74	1/3531 (0.0%)
1	C	0.40	0/2543	0.76	0/3472
2	B	0.38	0/2367	0.75	2/3223 (0.1%)
2	D	0.36	0/2404	0.73	2/3267 (0.1%)
All	All	0.38	0/9902	0.74	5/13493 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	96	GLU	N-CA-C	-5.38	105.08	111.69
2	D	240	VAL	CA-C-N	5.21	125.26	119.32
2	D	240	VAL	C-N-CA	5.21	125.26	119.32
2	B	240	VAL	CA-C-N	5.03	125.05	119.32
2	B	240	VAL	C-N-CA	5.03	125.05	119.32

There are no chirality outliers.

There are no planarity outliers.

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	2538	0	2552	63	0
1	C	2495	0	2448	62	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2332	0	2279	56	0
2	D	2369	0	2360	31	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	31	0	10	0	0
4	C	31	0	10	3	0
All	All	9798	0	9659	207	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 11.

All (207) 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:184:LEU:O	1:A:187:LEU:CD1	1.65	1.45
1:A:184:LEU:HA	1:A:187:LEU:HD11	1.36	1.03
1:A:184:LEU:O	1:A:187:LEU:HD13	1.56	1.03
2:B:221:ASN:ND2	2:B:356:THR:HG21	1.77	0.98
1:A:184:LEU:CA	1:A:187:LEU:HD11	1.92	0.98
1:A:184:LEU:O	1:A:187:LEU:HD11	1.62	0.95
1:A:184:LEU:C	1:A:187:LEU:HD11	1.91	0.94
1:A:184:LEU:C	1:A:187:LEU:CD1	2.40	0.94
1:A:184:LEU:O	1:A:187:LEU:HD12	1.64	0.94
1:C:89:ILE:HG23	1:C:99:ILE:O	1.68	0.92
1:A:94:THR:O	1:A:98:GLN:HG3	1.69	0.91
1:A:343:ILE:HD11	1:A:350:ILE:CG2	2.00	0.91
1:C:93:GLY:HA3	1:C:94:THR:CB	1.99	0.90
1:C:145:SER:O	1:C:149:GLN:HG3	1.74	0.87
1:A:359:GLU:O	1:A:360:LYS:HB2	1.77	0.85
1:C:93:GLY:CA	1:C:94:THR:CB	2.57	0.82
1:A:135:VAL:HG22	1:A:177:ILE:HD13	1.61	0.81
1:C:217:GLN:OE1	1:C:254:MET:HE1	1.82	0.79
2:D:310:GLN:O	2:D:314:ILE:HG13	1.81	0.79
1:A:186:LYS:CG	1:A:186:LYS:O	2.31	0.78
1:C:228:LYS:O	1:C:232:ILE:HG12	1.84	0.77
2:B:221:ASN:HD21	2:B:356:THR:HG21	1.48	0.77
1:C:217:GLN:OE1	1:C:254:MET:CE	2.33	0.76
1:C:135:VAL:HG22	1:C:177:ILE:HD13	1.66	0.76
1:A:45:LYS:HB2	1:A:46:PRO:HD2	1.68	0.75
2:D:182:LYS:HD2	2:D:189:LEU:HD23	1.69	0.75
2:B:182:LYS:NZ	2:B:182:LYS:HB2	2.03	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:53:GLU:O	2:B:53:GLU:HG2	1.89	0.72
2:B:221:ASN:CG	2:B:356:THR:HG21	2.16	0.70
1:C:92:PRO:O	1:C:93:GLY:C	2.35	0.70
2:B:373:PRO:HG2	1:C:254:MET:HE3	1.75	0.69
2:D:98:VAL:HG13	2:D:359:ILE:HD11	1.75	0.69
1:C:94:THR:O	1:C:95:PHE:HB2	1.91	0.68
2:B:17:LYS:HG2	2:B:334:LEU:HD12	1.77	0.67
2:D:236:LEU:O	2:D:236:LEU:HD12	1.94	0.67
1:C:256:ASN:OD1	1:C:258:THR:O	2.13	0.66
1:A:343:ILE:HD11	1:A:350:ILE:HG22	1.78	0.66
1:C:89:ILE:CG2	1:C:99:ILE:O	2.42	0.66
2:B:42:LEU:HD11	2:B:169:ILE:HD13	1.77	0.66
1:A:187:LEU:HD12	1:A:187:LEU:N	2.11	0.66
1:A:145:SER:O	1:A:149:GLN:HG3	1.97	0.64
1:A:87:ASN:O	1:A:91:ALA:HB2	1.97	0.64
1:C:231:ARG:HH11	1:C:232:ILE:CD1	2.11	0.64
1:A:110:LYS:HD3	1:A:211:PHE:CD2	2.34	0.63
1:A:186:LYS:O	1:A:186:LYS:HG2	1.97	0.63
1:C:78:VAL:O	1:C:82:ASN:HB2	1.99	0.63
1:C:231:ARG:NH1	1:C:232:ILE:CD1	2.62	0.62
2:B:7:PHE:HE2	2:B:270:GLU:HG3	1.64	0.62
1:C:231:ARG:NH1	1:C:232:ILE:HD13	2.14	0.62
2:B:7:PHE:CE2	2:B:270:GLU:HG3	2.34	0.62
1:C:231:ARG:HH11	1:C:232:ILE:HD13	1.66	0.61
1:C:45:LYS:HB2	1:C:46:PRO:HD2	1.81	0.61
1:C:29:HIS:HB3	1:C:283:GLY:O	2.01	0.61
2:B:182:LYS:HB2	2:B:182:LYS:HZ2	1.64	0.61
1:C:60:ARG:HD3	1:C:332:GLN:HG2	1.82	0.60
1:A:95:PHE:HA	1:A:98:GLN:HB2	1.81	0.60
2:D:92:VAL:O	2:D:93:MET:HE2	2.02	0.60
2:D:307:ASP:O	2:D:311:ARG:HG3	2.03	0.59
1:A:187:LEU:HD12	1:A:187:LEU:H	1.67	0.59
2:D:43:LYS:HA	2:D:365:TYR:OH	2.02	0.59
2:B:53:GLU:O	2:B:53:GLU:CG	2.51	0.59
1:A:29:HIS:HB3	1:A:283:GLY:O	2.03	0.59
2:D:187:GLU:O	2:D:187:GLU:HG3	2.03	0.58
1:C:91:ALA:O	1:C:93:GLY:N	2.30	0.58
2:B:367:ILE:O	2:B:369:PRO:HD3	2.04	0.57
2:B:181:GLU:O	2:B:185:ALA:HB3	2.04	0.57
1:C:31:PRO:HG2	1:C:279:LEU:HA	1.88	0.56
2:D:42:LEU:HD11	2:D:169:ILE:HD13	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:309:MET:HA	1:C:312:TYR:CE2	2.42	0.55
2:D:124:LEU:O	2:D:128:VAL:HG23	2.06	0.55
2:D:182:LYS:CD	2:D:189:LEU:HD23	2.38	0.54
2:D:49:ILE:O	2:D:53:GLU:HG3	2.08	0.54
1:C:99:ILE:HG23	1:C:124:VAL:CG1	2.38	0.54
1:A:160:ASN:OD1	1:A:163:GLY:C	2.50	0.54
2:D:235:ASP:O	2:D:239:ARG:HG2	2.08	0.54
2:B:131:ASN:O	2:B:135:GLN:HG2	2.08	0.54
2:B:330:LEU:HG	2:B:334:LEU:HD23	1.90	0.53
1:A:217:GLN:HG2	1:A:220:VAL:HG23	1.89	0.53
1:A:309:MET:HA	1:A:312:TYR:CE2	2.44	0.53
2:B:132:LEU:HD21	2:B:157:ILE:HD11	1.89	0.53
2:D:107:ASP:OD2	2:D:234:ARG:NH2	2.41	0.53
1:A:90:VAL:O	1:A:92:PRO:HD3	2.09	0.53
2:B:186:GLY:O	2:B:187:GLU:C	2.52	0.53
1:C:287:ILE:O	1:C:298:VAL:HG22	2.09	0.52
1:A:87:ASN:O	1:A:91:ALA:CB	2.57	0.52
1:C:217:GLN:OE1	1:C:254:MET:HE2	2.10	0.52
1:A:186:LYS:O	1:A:186:LYS:HG3	2.07	0.52
1:A:228:LYS:O	1:A:232:ILE:HG12	2.10	0.51
2:B:221:ASN:HD21	2:B:356:THR:CG2	2.22	0.51
1:C:60:ARG:HD3	1:C:332:GLN:CG	2.40	0.51
2:B:221:ASN:CG	2:B:356:THR:CG2	2.84	0.51
2:B:221:ASN:OD1	2:B:356:THR:CG2	2.59	0.51
1:C:316:THR:O	1:C:320:ILE:HG12	2.12	0.50
1:A:287:ILE:O	1:A:298:VAL:HG22	2.11	0.50
2:B:7:PHE:HD2	2:B:270:GLU:HB2	1.76	0.50
2:D:48:TRP:O	2:D:52:GLN:HB2	2.12	0.50
1:C:254:MET:HE1	1:C:261:PRO:HG3	1.94	0.49
2:B:181:GLU:O	2:B:185:ALA:CB	2.60	0.49
2:B:183:VAL:C	2:B:186:GLY:H	2.20	0.49
1:A:178:THR:HA	1:A:198:LEU:HD22	1.95	0.49
1:C:42:PRO:HG2	2:D:292:TYR:CB	2.42	0.49
1:C:100:GLU:OE1	1:C:127:LYS:HE3	2.13	0.49
1:A:220:VAL:O	1:A:224:ILE:HG13	2.13	0.49
1:C:85:ILE:O	1:C:89:ILE:CG1	2.61	0.48
2:B:182:LYS:O	2:B:183:VAL:C	2.55	0.48
1:C:86:ASP:O	1:C:90:VAL:HG23	2.13	0.48
2:B:6:ILE:O	2:B:7:PHE:HB2	2.14	0.48
2:B:225:SER:O	2:B:229:VAL:HG23	2.13	0.48
1:A:60:ARG:HD3	1:A:332:GLN:HG2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:319:ARG:HD2	2:D:312:GLU:OE2	2.13	0.48
1:A:258:THR:O	1:A:259:ARG:CB	2.61	0.48
2:B:221:ASN:ND2	2:B:356:THR:CG2	2.65	0.48
2:B:11:ASP:O	2:B:15:MET:HG2	2.14	0.47
2:B:45:VAL:CG2	2:B:135:GLN:HG3	2.44	0.47
1:C:57:LEU:HD23	1:C:330:LEU:HD21	1.96	0.47
2:D:103:LEU:O	2:D:104:LEU:HD23	2.13	0.47
2:B:235:ASP:O	2:B:239:ARG:HG2	2.15	0.47
2:B:249:TRP:HB3	2:B:250:PRO:HD3	1.97	0.47
1:C:233:ARG:NH2	1:C:329:ILE:O	2.38	0.47
1:C:76:SER:O	1:C:80:LYS:HG3	2.13	0.47
1:C:178:THR:HB	1:C:201:ALA:HB3	1.97	0.46
1:A:343:ILE:HD11	1:A:350:ILE:HG21	1.91	0.46
1:C:210:TRP:CD1	1:C:214:ASN:HD22	2.33	0.46
1:A:308:ASP:OD1	2:B:323:ARG:HD2	2.15	0.46
1:A:31:PRO:HG2	1:A:279:LEU:HA	1.97	0.46
1:A:129:LEU:HD13	1:A:192:HIS:CE1	2.51	0.46
2:B:179:GLU:O	2:B:180:MET:C	2.59	0.46
2:B:182:LYS:NZ	2:B:182:LYS:CB	2.73	0.46
2:B:226:CYS:O	2:B:230:ILE:HG13	2.16	0.46
1:C:85:ILE:HG23	1:C:99:ILE:HD13	1.98	0.46
2:B:307:ASP:O	2:B:311:ARG:HG3	2.16	0.46
1:A:39:MET:HE2	1:C:38:GLU:OE2	2.16	0.45
1:C:220:VAL:O	1:C:224:ILE:HG13	2.16	0.45
2:D:249:TRP:HB3	2:D:250:PRO:HD3	1.98	0.45
1:A:178:THR:HB	1:A:201:ALA:HB3	1.97	0.45
1:A:336:ALA:O	1:A:339:LEU:HB2	2.17	0.45
1:A:167:SER:OG	1:A:172:THR:OG1	2.26	0.45
2:D:196:ASP:OD1	2:D:196:ASP:N	2.50	0.45
2:B:198:LEU:HD12	2:B:198:LEU:HA	1.83	0.45
1:A:167:SER:OG	1:A:172:THR:CG2	2.65	0.44
2:D:187:GLU:O	2:D:187:GLU:CG	2.65	0.44
1:C:100:GLU:HB2	1:C:127:LYS:HG2	1.99	0.44
1:C:308:ASP:OD1	2:D:323:ARG:HD2	2.17	0.44
1:A:238:GLU:HB3	1:A:239:PRO:HD3	2.00	0.44
1:C:238:GLU:HB3	1:C:239:PRO:HD3	1.99	0.44
2:D:167:LEU:C	2:D:167:LEU:HD23	2.42	0.44
1:A:181:PRO:HB2	1:A:182:PRO:HD3	2.00	0.44
2:B:167:LEU:C	2:B:167:LEU:HD23	2.41	0.44
1:A:129:LEU:HB3	1:A:192:HIS:CD2	2.52	0.44
2:B:373:PRO:HG2	1:C:254:MET:CE	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:MET:HE3	1:C:157:MET:HB3	1.85	0.44
1:A:45:LYS:HB2	1:A:46:PRO:CD	2.41	0.43
2:B:45:VAL:HG23	2:B:135:GLN:HG3	1.99	0.43
2:B:103:LEU:O	2:B:104:LEU:HD23	2.18	0.43
1:A:272:LEU:HD12	1:A:321:LEU:HD22	1.99	0.43
2:B:40:ARG:HG3	2:B:364:THR:OG1	2.18	0.43
1:A:110:LYS:HD3	1:A:211:PHE:CG	2.54	0.43
2:B:257:LYS:O	2:B:261:THR:HG23	2.17	0.43
1:C:80:LYS:O	1:C:84:VAL:HG23	2.18	0.43
2:D:53:GLU:C	2:D:55:GLY:N	2.77	0.43
2:D:103:LEU:HB2	2:D:358:GLN:OE1	2.19	0.43
1:A:251:TYR:C	1:A:251:TYR:CD1	2.97	0.43
1:C:98:GLN:HB3	1:C:127:LYS:HB2	2.00	0.43
1:C:104:GLN:HG2	1:C:122:LEU:HD21	1.99	0.43
4:C:1362:ATP:O1B	4:C:1362:ATP:O1A	2.37	0.43
2:D:11:ASP:O	2:D:15:MET:HG2	2.19	0.43
1:A:104:GLN:HG2	1:A:122:LEU:HD21	2.01	0.42
2:B:162:GLU:HA	2:B:163:PRO:HA	1.85	0.42
1:C:135:VAL:HG13	1:C:177:ILE:CD1	2.49	0.42
1:C:293:SER:HA	1:C:294:GLY:HA2	1.66	0.42
1:A:91:ALA:O	1:A:92:PRO:C	2.61	0.42
2:B:310:GLN:O	2:B:314:ILE:HG13	2.19	0.42
1:A:162:THR:HA	1:A:197:VAL:HG11	2.02	0.42
4:C:1362:ATP:O1B	4:C:1362:ATP:O1G	2.38	0.42
2:D:120:PRO:HB2	2:D:198:LEU:HD12	2.02	0.42
2:D:22:TYR:HA	2:D:23:PRO:HD3	1.89	0.42
1:A:95:PHE:CE2	1:A:141:LYS:HG2	2.55	0.41
1:C:90:VAL:O	1:C:90:VAL:HG12	2.20	0.41
1:C:231:ARG:HH11	1:C:232:ILE:HD11	1.84	0.41
4:C:1362:ATP:H8	4:C:1362:ATP:O5'	2.03	0.41
2:D:157:ILE:HD11	2:D:169:ILE:HD11	2.02	0.41
1:A:349:VAL:HG21	2:B:308:ARG:HB3	2.03	0.41
2:B:34:MET:HE3	2:B:157:ILE:HG22	2.03	0.41
2:B:372:ARG:HA	2:B:373:PRO:HD3	1.90	0.41
1:C:98:GLN:CB	1:C:127:LYS:HB2	2.50	0.41
1:A:135:VAL:HG13	1:A:177:ILE:CD1	2.49	0.41
1:A:319:ARG:NH2	1:A:354:SER:OG	2.53	0.41
1:A:343:ILE:O	1:A:343:ILE:HG23	2.20	0.41
1:C:180:VAL:HB	1:C:182:PRO:HD2	2.02	0.41
2:B:372:ARG:HG2	2:B:373:PRO:HD2	2.03	0.41
1:C:85:ILE:O	1:C:89:ILE:HG12	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:224:LYS:HD3	2:B:266:MET:O	2.20	0.41
1:C:75:LEU:O	1:C:79:THR:OG1	2.31	0.41
1:C:233:ARG:NH1	1:C:233:ARG:HB3	2.36	0.41
2:B:285:MET:HE3	2:B:286:PRO:HD2	2.02	0.41
1:C:278:GLY:HA2	1:C:281:LEU:HG	2.03	0.41
1:A:311:CYS:O	1:A:315:GLN:HG3	2.21	0.40
2:B:266:MET:HE2	2:B:270:GLU:OE2	2.19	0.40
2:D:215:TRP:CH2	2:D:253:LEU:HD23	2.56	0.40
1:C:60:ARG:HG3	1:C:330:LEU:HB3	2.02	0.40
2:D:162:GLU:HA	2:D:163:PRO:HA	1.89	0.40
1:A:188:ASP:O	1:A:190:GLU:N	2.55	0.40
1:A:278:GLY:HA2	1:A:281:LEU:HG	2.01	0.40
2:B:177:ARG:NH1	2:B:181:GLU:OE2	2.53	0.40
1:A:356:LYS:HE2	1:A:358:TYR:O	2.21	0.40
2:B:224:LYS:O	2:B:225:SER:HB2	2.21	0.40
2:B:248:GLY:O	2:B:252:GLU:HG3	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	331/364 (91%)	306 (92%)	24 (7%)	1 (0%)	36	60
1	C	328/364 (90%)	305 (93%)	20 (6%)	3 (1%)	14	35
2	B	304/383 (79%)	293 (96%)	11 (4%)	0	100	100
2	D	304/383 (79%)	292 (96%)	12 (4%)	0	100	100
All	All	1267/1494 (85%)	1196 (94%)	67 (5%)	4 (0%)	36	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	257	PRO
1	C	92	PRO
1	A	92	PRO
1	C	93	GLY

5.3.2 Protein sidechains [i](#)

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	274/317 (86%)	269 (98%)	5 (2%)	51	78
1	C	262/317 (83%)	253 (97%)	9 (3%)	32	62
2	B	239/330 (72%)	233 (98%)	6 (2%)	42	71
2	D	249/330 (76%)	242 (97%)	7 (3%)	38	68
All	All	1024/1294 (79%)	997 (97%)	27 (3%)	40	70

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	ASN
1	A	177	ILE
1	A	217	GLN
1	A	339	LEU
1	A	350	ILE
2	B	53	GLU
2	B	93	MET
2	B	111	GLU
2	B	182	LYS
2	B	198	LEU
2	B	214	LYS
1	C	82	ASN
1	C	89	ILE
1	C	160	ASN
1	C	161	GLU
1	C	177	ILE
1	C	228	LYS
1	C	343	ILE

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Mol	Chain	Res	Type
1	C	345	THR
1	C	356	LYS
2	D	111	GLU
2	D	127	LYS
2	D	157	ILE
2	D	189	LEU
2	D	236	LEU
2	D	314	ILE
2	D	357	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	250	HIS
1	A	299	HIS
2	B	212	HIS
1	C	140	ASN
1	C	250	HIS
1	C	332	GLN
2	D	131	ASN
2	D	135	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 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	$\# Z > 2$
4	ATP	C	1362	3	32,33,33	1.83	8 (25%)	48,52,52	1.90	10 (20%)
4	ATP	A	1362	3	32,33,33	1.73	8 (25%)	48,52,52	1.97	12 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ATP	C	1362	3	-	4/22/38/38	0/3/3/3
4	ATP	A	1362	3	-	5/22/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1362	ATP	PB-O3A	-5.46	1.53	1.59
4	A	1362	ATP	PB-O3A	-4.55	1.54	1.59
4	C	1362	ATP	C6-N6	3.61	1.43	1.34
4	A	1362	ATP	C6-N6	3.57	1.43	1.34
4	A	1362	ATP	O2'-C2'	-3.45	1.34	1.43
4	C	1362	ATP	O2'-C2'	-3.42	1.34	1.43
4	C	1362	ATP	O3'-C3'	-3.11	1.35	1.43
4	A	1362	ATP	O3'-C3'	-3.04	1.35	1.43
4	A	1362	ATP	PB-O3B	-2.75	1.56	1.59
4	C	1362	ATP	PB-O3B	-2.74	1.56	1.59
4	C	1362	ATP	C2'-C3'	-2.59	1.46	1.53
4	A	1362	ATP	C5'-C4'	-2.27	1.44	1.51
4	A	1362	ATP	C2'-C3'	-2.24	1.47	1.53
4	A	1362	ATP	C3'-C4'	-2.10	1.47	1.53
4	C	1362	ATP	C5'-C4'	-2.08	1.45	1.51
4	C	1362	ATP	C3'-C4'	-2.04	1.47	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1362	ATP	C5-C4-N3	-5.36	119.33	126.72
4	C	1362	ATP	C5-C4-N3	-5.24	119.50	126.72
4	A	1362	ATP	N3-C2-N1	-4.86	121.23	128.58
4	C	1362	ATP	N3-C2-N1	-4.75	121.39	128.58
4	C	1362	ATP	N3-C4-N9	4.49	134.80	127.17
4	A	1362	ATP	N3-C4-N9	4.27	134.43	127.17
4	C	1362	ATP	N9-C8-N7	-4.16	108.03	113.94
4	A	1362	ATP	N9-C8-N7	-4.05	108.19	113.94
4	C	1362	ATP	C4-N9-C8	3.80	109.73	105.74
4	A	1362	ATP	C2-N3-C4	3.59	120.60	111.83
4	A	1362	ATP	C5-N7-C8	3.59	109.09	103.45
4	C	1362	ATP	C5-N7-C8	3.50	108.95	103.45
4	C	1362	ATP	C2-N3-C4	3.48	120.33	111.83
4	A	1362	ATP	C4-N9-C8	3.35	109.26	105.74
4	A	1362	ATP	C4-C5-N7	-2.93	107.23	110.58
4	C	1362	ATP	O5'-C5'-C4'	2.87	118.75	108.99
4	A	1362	ATP	C3'-C2'-C1'	2.81	106.78	101.46
4	C	1362	ATP	C4-C5-N7	-2.62	107.58	110.58
4	C	1362	ATP	O3G-PG-O3B	2.48	112.94	104.64
4	A	1362	ATP	C2'-C3'-C4'	2.45	107.33	102.61
4	A	1362	ATP	O3G-PG-O3B	2.24	112.14	104.64
4	A	1362	ATP	O5'-C5'-C4'	2.14	116.29	108.99

There are no chirality outliers.

All (9) torsion outliers are listed below:

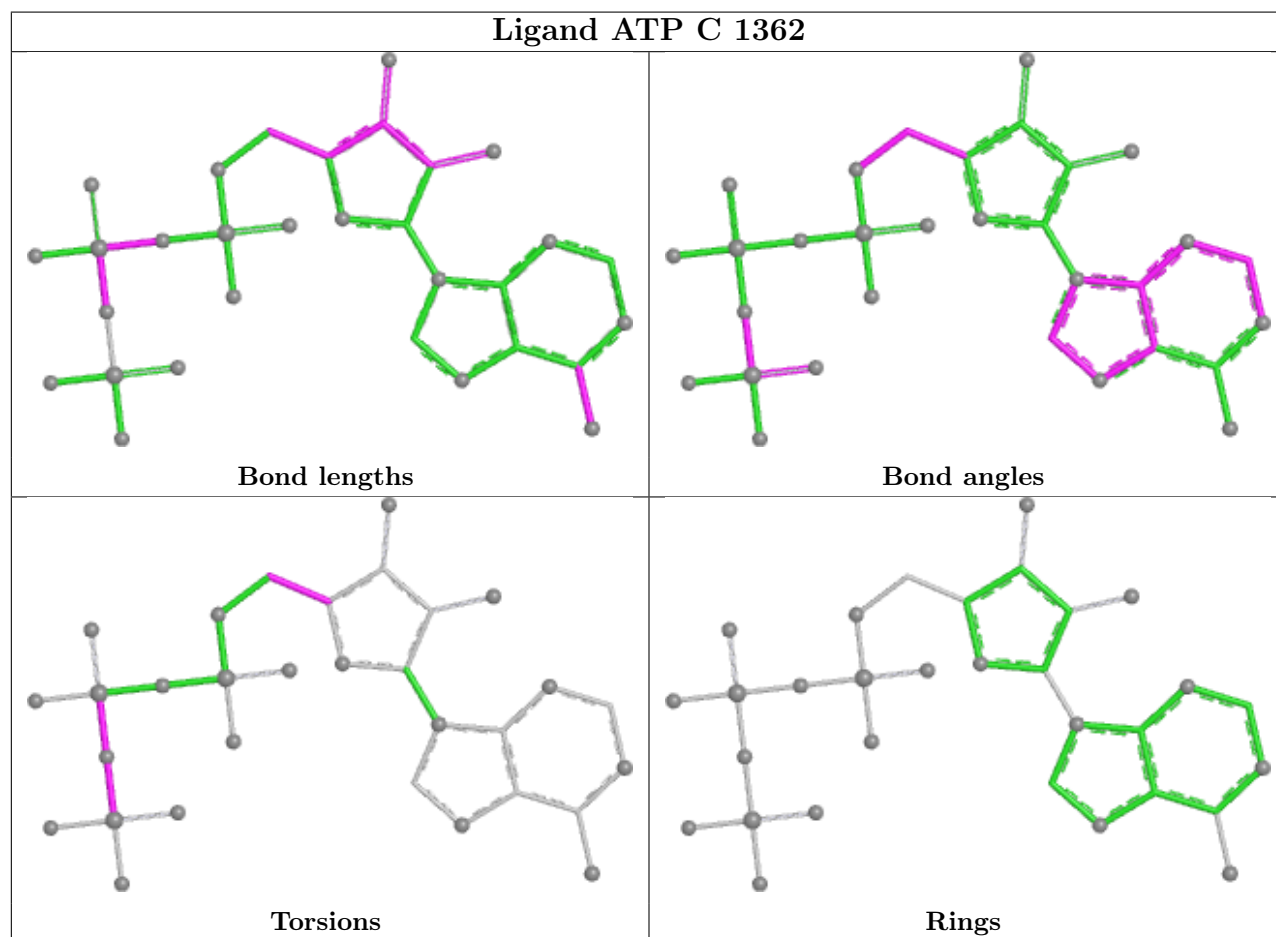
Mol	Chain	Res	Type	Atoms
4	A	1362	ATP	O4'-C4'-C5'-O5'
4	A	1362	ATP	C3'-C4'-C5'-O5'
4	C	1362	ATP	PB-O3B-PG-O3G
4	A	1362	ATP	PB-O3A-PA-O1A
4	C	1362	ATP	PB-O3B-PG-O1G
4	C	1362	ATP	O4'-C4'-C5'-O5'
4	A	1362	ATP	PG-O3B-PB-O1B
4	C	1362	ATP	PG-O3B-PB-O2B
4	A	1362	ATP	PB-O3A-PA-O2A

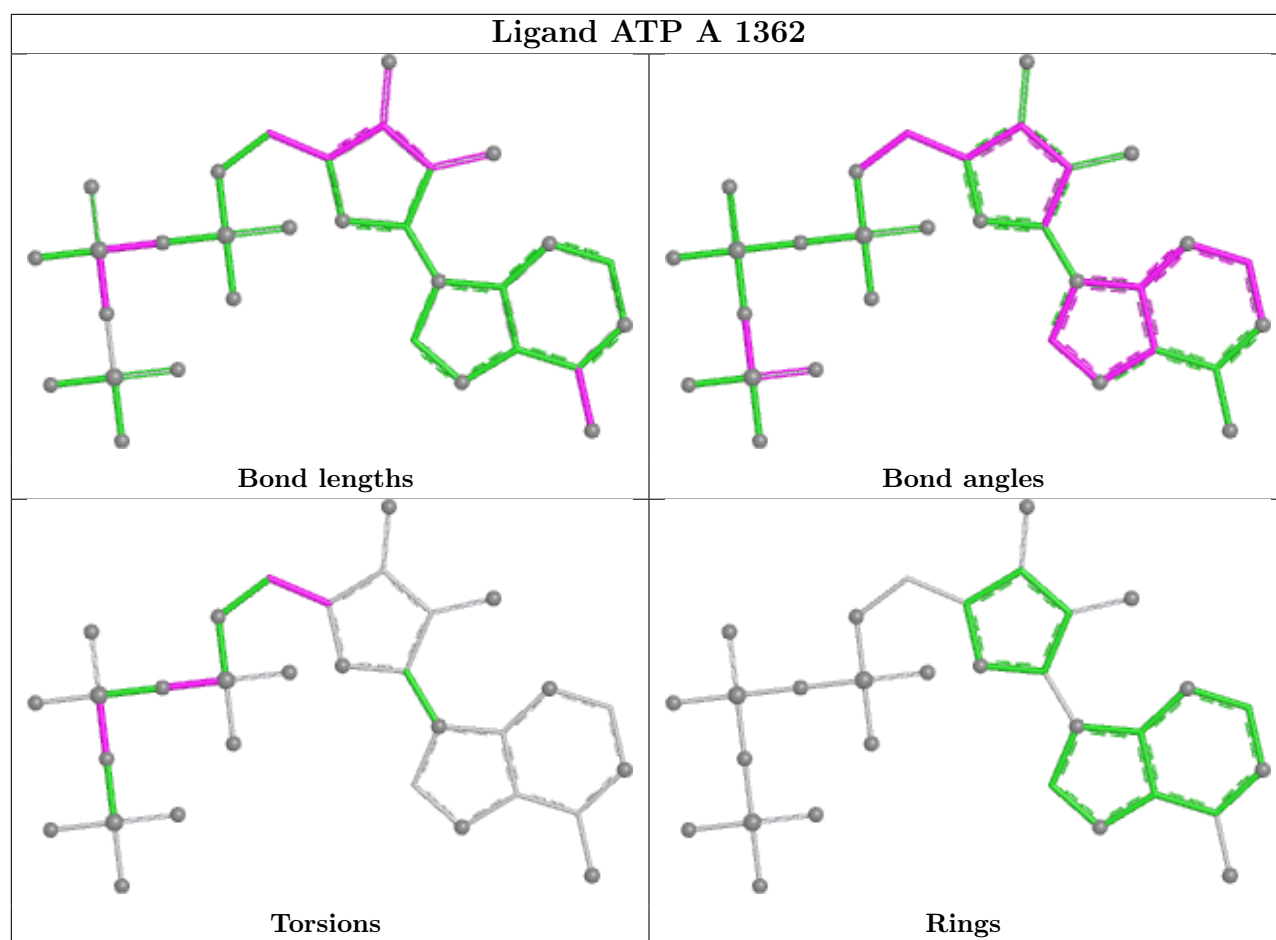
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1362	ATP	3	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	333/364 (91%)	0.45	28 (8%) 17 14	11, 28, 65, 116	0
1	C	332/364 (91%)	0.69	34 (10%) 12 10	13, 35, 79, 123	0
2	B	316/383 (82%)	0.65	39 (12%) 8 7	11, 35, 85, 125	0
2	D	316/383 (82%)	0.46	24 (7%) 20 17	12, 31, 73, 114	0
All	All	1297/1494 (86%)	0.56	125 (9%) 13 11	11, 32, 75, 125	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	97	VAL	7.2
2	B	373	PRO	6.8
2	D	373	PRO	6.7
2	B	354	ASP	5.9
2	D	370	MET	5.9
1	A	260	GLN	5.8
2	D	189	LEU	5.5
2	B	355	TYR	5.4
2	B	369	PRO	5.3
2	B	368	THR	5.3
2	B	367	ILE	5.2
2	D	369	PRO	5.1
2	D	263	ASN	5.1
2	D	195	PRO	5.0
2	D	6	ILE	5.0
1	A	92	PRO	4.9
2	D	372	ARG	4.9
1	C	258	THR	4.8
1	C	217	GLN	4.8
1	C	361	PRO	4.7
1	A	360	LYS	4.6

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Mol	Chain	Res	Type	RSRZ
2	B	195	PRO	4.5
2	D	366	ALA	4.5
2	B	372	ARG	4.4
2	D	87	ARG	4.3
2	D	162	GLU	4.2
1	C	216	SER	4.2
1	A	68	SER	4.1
2	D	188	THR	4.1
1	C	129	LEU	4.1
2	B	87	ARG	4.1
2	B	109	ASP	4.1
1	C	94	THR	4.0
2	B	363	THR	4.0
2	D	340	PRO	4.0
1	C	259	ARG	3.9
1	C	260	GLN	3.9
2	B	162	GLU	3.9
1	A	93	GLY	3.8
2	B	365	TYR	3.7
1	C	180	VAL	3.7
1	A	217	GLN	3.6
1	A	172	THR	3.6
2	B	186	GLY	3.6
1	A	293	SER	3.6
2	B	6	ILE	3.6
1	C	43	ARG	3.6
2	B	161	LYS	3.5
2	D	55	GLY	3.5
2	D	365	TYR	3.4
2	D	354	ASP	3.3
1	C	68	SER	3.2
1	A	91	ALA	3.2
2	B	183	VAL	3.2
2	B	193	ASP	3.1
1	C	257	PRO	3.1
1	A	161	GLU	3.1
1	A	258	THR	3.0
2	B	366	ALA	3.0
1	A	185	ARG	3.0
2	D	357	VAL	2.9
1	C	92	PRO	2.9
2	B	357	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	153	GLU	2.9
2	B	54	LYS	2.9
2	D	150	VAL	2.9
2	B	187	GLU	2.8
1	A	294	GLY	2.8
1	C	182	PRO	2.8
1	A	216	SER	2.8
2	B	197	VAL	2.8
1	C	360	LYS	2.8
2	B	108	LEU	2.8
2	D	50	ASP	2.8
1	C	181	PRO	2.7
1	C	128	ILE	2.7
1	C	293	SER	2.7
2	B	50	ASP	2.7
2	B	198	LEU	2.7
1	C	39	MET	2.7
1	C	93	GLY	2.7
1	C	161	GLU	2.7
2	B	182	LYS	2.7
2	B	177	ARG	2.7
1	A	187	LEU	2.6
1	A	0	SER	2.6
1	C	99	ILE	2.6
1	C	136	ALA	2.6
2	B	150	VAL	2.6
2	B	364	THR	2.5
1	A	44	VAL	2.5
1	C	255	ASN	2.5
2	D	166	SER	2.5
1	C	295	ASN	2.5
1	A	259	ARG	2.4
1	C	185	ARG	2.4
2	B	194	PRO	2.4
2	D	355	TYR	2.4
2	B	53	GLU	2.4
2	B	298	GLU	2.4
2	D	187	GLU	2.4
1	A	73	SER	2.3
2	B	180	MET	2.3
2	B	340	PRO	2.3
1	C	90	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	190	GLU	2.3
1	C	89	ILE	2.3
1	A	181	PRO	2.3
2	B	92	VAL	2.3
1	A	47	ALA	2.2
1	C	170	ASP	2.2
1	A	46	PRO	2.2
2	D	108	LEU	2.2
1	A	95	PHE	2.2
2	B	356	THR	2.2
2	D	287	ASP	2.2
1	A	213	GLU	2.1
1	C	213	GLU	2.1
1	A	343	ILE	2.1
2	B	21	VAL	2.1
2	B	7	PHE	2.1
1	C	107	SER	2.0
1	A	345	THR	2.0
1	C	130	PRO	2.0
1	C	67	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	1361	1/1	0.81	0.23	51,51,51,51	0
3	MG	C	1363	1/1	0.81	0.20	51,51,51,51	0
4	ATP	C	1362	31/31	0.88	0.16	41,41,41,41	31

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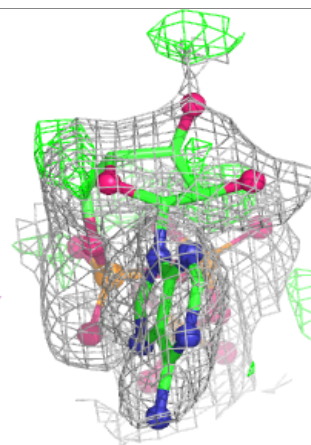
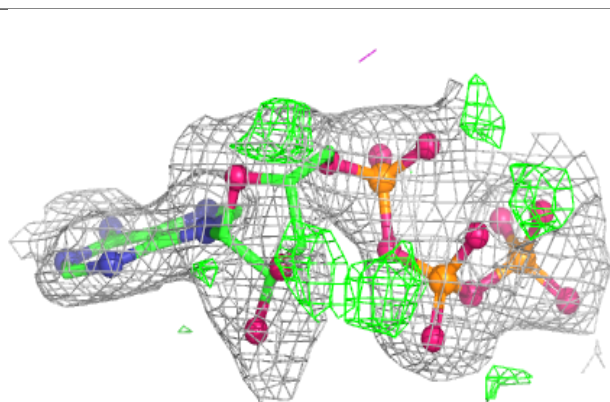
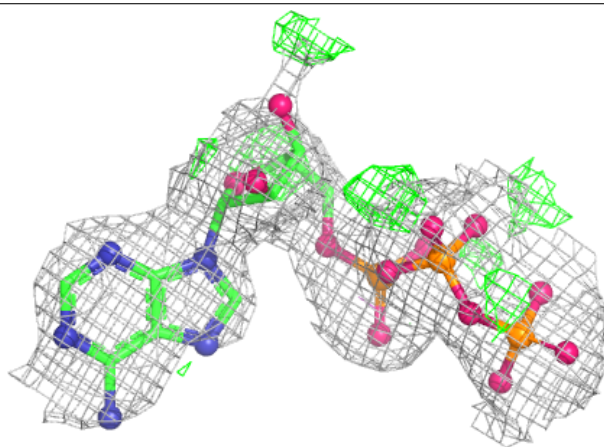
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ATP	A	1362	31/31	0.90	0.14	41,41,41,41	31

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.

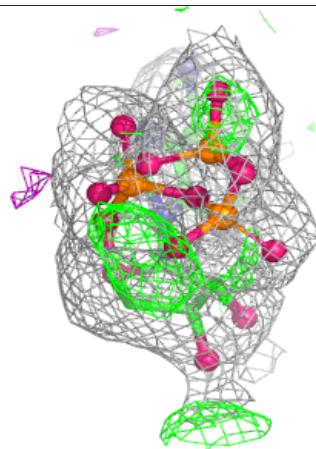
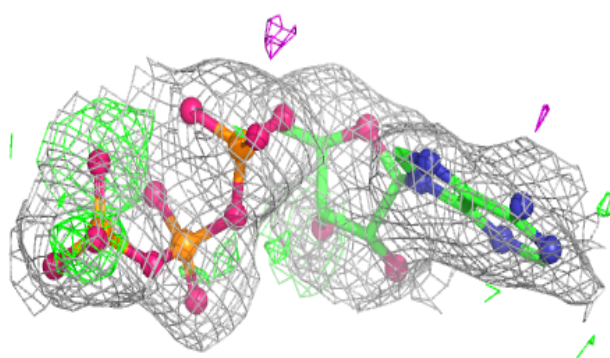
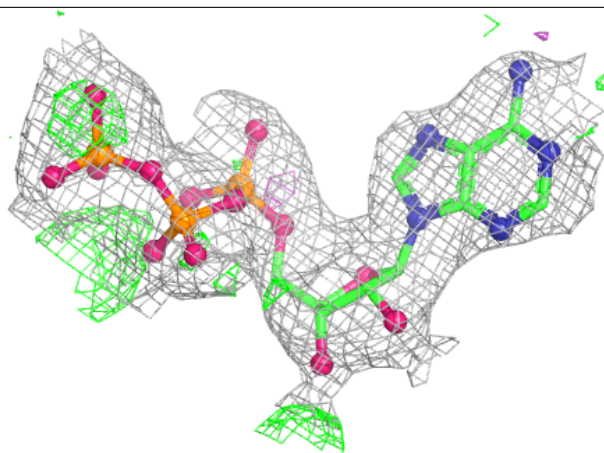
Electron density around ATP C 1362:

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 1362:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.