Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2021 – 10:38 AM EDT

PDB ID	:	1GOL
Title	:	COORDINATES OF RAT MAP KINASE ERK2 WITH AN ARGININE MUTATION AT POSITION 52
Authors	:	Harkins, P.C.; Zhang, F.; Goldsmith, E.J.
Deposited on	:	1996-01-23
Resolution	:	2.80 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 (i) symbol.

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

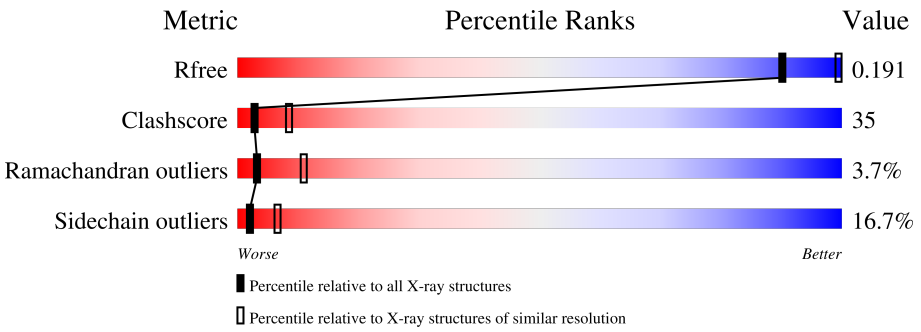
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	364	49% 37% 10% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3012 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 EXTRACELLULAR REGULATED KINASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	357	2901	1859	499	528	15	0	0	0

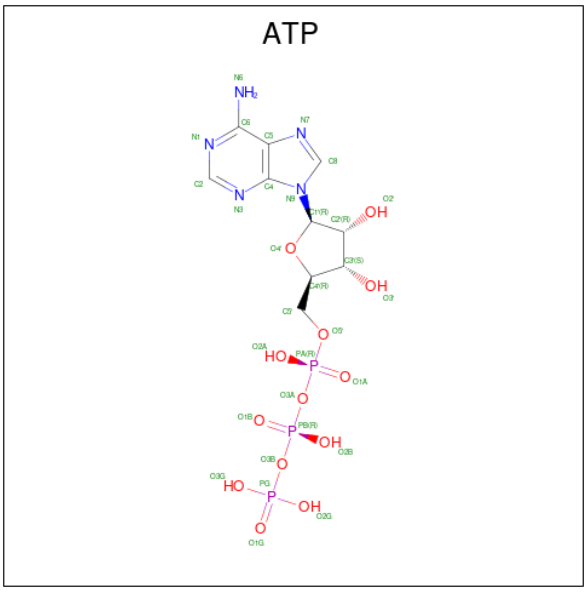
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	ARG	LYS	engineered mutation	UNP P63086

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	1	1	1	0	0

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 4 is water.

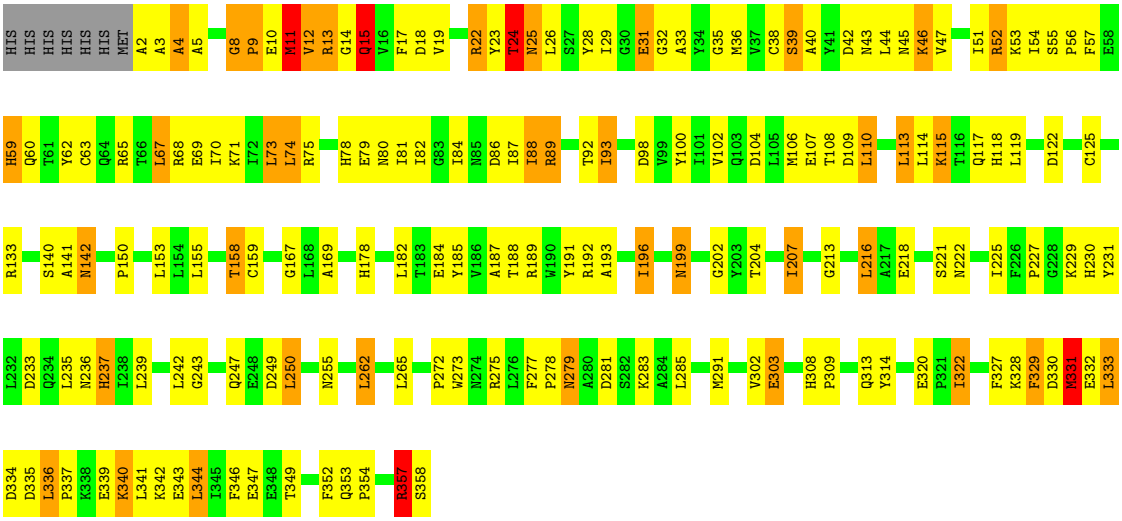
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	79	79	79	0	0

3 Residue-property plots (i)

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. 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. 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: EXTRACELLULAR REGULATED KINASE 2

Chain A: 49% 37% 10%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	49.32Å 71.42Å 61.25Å 90.00° 109.75° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.80 15.69 – 1.81	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.80) 75.3 (15.69-1.81)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.07 (at 1.81Å)	Xtriage
Refinement program		Depositor
R, R_{free}	(Not available) , (Not available) 0.189 , 0.191	Depositor DCC
R_{free} test set	1395 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	19.8	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 101.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3012	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.14% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, 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	0.49	1/2971 (0.0%)	0.65	1/4026 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	8	GLY	C-N	12.95	1.58	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	52	ARG	NE-CZ-NH2	-5.27	117.67	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	191	TYR	Sidechain

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	2901	0	2897	202	0
2	A	1	0	0	0	0
3	A	31	0	12	7	0
4	A	79	0	0	16	0
All	All	3012	0	2909	206	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 35.

All (206) 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)
3:A:581:ATP:C8	3:A:581:ATP:N7	1.94	1.35
3:A:581:ATP:O4'	3:A:581:ATP:C1'	1.65	1.32
1:A:207:ILE:HD13	4:A:514:HOH:O	1.22	1.31
1:A:14:GLY:HA2	1:A:36:MET:SD	1.80	1.19
1:A:12:VAL:HG12	1:A:13:ARG:HE	1.21	1.02
1:A:9:PRO:HA	1:A:14:GLY:HA3	1.45	0.99
1:A:13:ARG:HA	1:A:15:GLN:HE22	1.27	0.98
1:A:12:VAL:HG12	1:A:13:ARG:NE	1.79	0.97
1:A:79:GLU:HG3	4:A:539:HOH:O	1.68	0.91
1:A:327:PHE:O	1:A:328:LYS:HG3	1.72	0.90
1:A:9:PRO:HA	1:A:14:GLY:CA	2.03	0.88
1:A:336:LEU:HD22	1:A:337:PRO:HD2	1.55	0.88
1:A:14:GLY:CA	1:A:36:MET:SD	2.63	0.86
1:A:78:HIS:HD2	1:A:80:ASN:H	1.19	0.84
1:A:207:ILE:CD1	4:A:514:HOH:O	1.96	0.83
1:A:185:TYR:HE2	1:A:187:ALA:HB3	1.42	0.83
1:A:185:TYR:CE2	1:A:187:ALA:HB3	2.15	0.82
1:A:340:LYS:NZ	1:A:344:LEU:HD13	1.95	0.81
1:A:52:ARG:HH11	1:A:52:ARG:HG3	1.47	0.79
1:A:340:LYS:HZ2	1:A:344:LEU:HD13	1.43	0.78
1:A:12:VAL:HG11	1:A:32:GLY:O	1.83	0.78
1:A:3:ALA:C	1:A:5:ALA:H	1.86	0.77
1:A:93:ILE:HD11	1:A:343:GLU:HA	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:ASN:O	4:A:813:HOH:O	2.03	0.76
1:A:14:GLY:HA2	1:A:36:MET:CE	2.15	0.75
1:A:51:ILE:HG12	1:A:102:VAL:HG22	1.68	0.74
1:A:22:ARG:HG2	1:A:23:TYR:CZ	2.22	0.74
1:A:57:PHE:CD1	1:A:342:LYS:HG3	2.23	0.73
1:A:331:MET:O	1:A:333:LEU:N	2.21	0.73
1:A:113:LEU:HD22	1:A:117:GLN:HB2	1.71	0.73
1:A:13:ARG:HB2	1:A:53:LYS:NZ	2.03	0.73
1:A:19:VAL:CG2	1:A:23:TYR:HB2	2.19	0.73
1:A:15:GLN:H	1:A:15:GLN:CD	1.92	0.72
1:A:330:ASP:O	1:A:331:MET:HB3	1.87	0.72
1:A:9:PRO:HG3	1:A:15:GLN:HG3	1.71	0.71
1:A:9:PRO:HB3	1:A:15:GLN:CD	2.11	0.70
1:A:193:ALA:HB3	1:A:196:ILE:HD12	1.73	0.70
1:A:13:ARG:CA	1:A:15:GLN:HE22	2.02	0.70
1:A:9:PRO:CB	1:A:15:GLN:HG3	2.21	0.70
1:A:57:PHE:HA	1:A:63:CYS:SG	2.32	0.70
1:A:15:GLN:H	1:A:15:GLN:NE2	1.89	0.69
1:A:140:SER:CB	1:A:322:ILE:HG12	2.22	0.69
1:A:78:HIS:CD2	1:A:80:ASN:H	2.08	0.69
1:A:9:PRO:HB3	1:A:15:GLN:CG	2.23	0.68
1:A:140:SER:HB2	1:A:322:ILE:HG12	1.76	0.67
1:A:188:THR:HB	4:A:513:HOH:O	1.95	0.66
1:A:26:LEU:HD23	1:A:39:SER:O	1.94	0.66
1:A:2:ALA:C	1:A:4:ALA:H	1.98	0.66
1:A:52:ARG:HG3	1:A:52:ARG:NH1	2.11	0.66
1:A:25:ASN:O	1:A:40:ALA:HA	1.95	0.66
1:A:24:THR:OG1	1:A:43:ASN:ND2	2.29	0.66
1:A:336:LEU:HD12	1:A:341:LEU:CD2	2.27	0.65
1:A:227:PRO:HD2	1:A:237:HIS:CD2	2.32	0.65
1:A:19:VAL:HG21	1:A:23:TYR:HB2	1.79	0.65
1:A:192:ARG:NH1	4:A:610:HOH:O	2.30	0.64
1:A:207:ILE:HD12	4:A:546:HOH:O	1.96	0.64
1:A:9:PRO:HB3	1:A:15:GLN:N	2.13	0.64
1:A:13:ARG:HA	1:A:15:GLN:NE2	2.09	0.64
1:A:17:PHE:HD2	1:A:36:MET:CE	2.11	0.64
1:A:45:ASN:O	1:A:47:VAL:HG13	1.99	0.63
1:A:357:ARG:HD3	1:A:358:SER:OG	1.98	0.63
1:A:9:PRO:HB3	1:A:15:GLN:HG3	1.80	0.63
1:A:9:PRO:HB3	1:A:15:GLN:OE1	1.98	0.62
1:A:9:PRO:CG	1:A:15:GLN:HG3	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:LYS:NZ	1:A:328:LYS:CB	2.63	0.62
1:A:17:PHE:CD2	1:A:36:MET:CE	2.83	0.62
1:A:22:ARG:O	1:A:22:ARG:HG3	2.00	0.61
1:A:43:ASN:O	1:A:46:LYS:HG2	2.00	0.61
1:A:42:ASP:O	1:A:46:LYS:N	2.33	0.61
1:A:9:PRO:HD3	1:A:15:GLN:HA	1.83	0.60
1:A:17:PHE:HD2	1:A:36:MET:HE2	1.65	0.60
1:A:322:ILE:CD1	4:A:709:HOH:O	2.49	0.60
1:A:13:ARG:HD2	1:A:35:GLY:HA2	1.83	0.60
1:A:19:VAL:HG23	1:A:23:TYR:HB2	1.84	0.60
1:A:13:ARG:HB2	1:A:53:LYS:HZ1	1.67	0.59
1:A:3:ALA:C	1:A:5:ALA:N	2.56	0.59
1:A:54:ILE:HG22	1:A:56:PRO:HD3	1.85	0.59
1:A:204:THR:OG1	4:A:514:HOH:O	2.08	0.59
1:A:28:TYR:HA	1:A:38:CYS:SG	2.43	0.59
1:A:15:GLN:CD	1:A:15:GLN:N	2.56	0.58
1:A:78:HIS:HB3	1:A:81:ILE:HD12	1.85	0.58
1:A:24:THR:O	1:A:25:ASN:C	2.41	0.58
1:A:340:LYS:CE	1:A:344:LEU:HD13	2.34	0.58
1:A:337:PRO:C	1:A:339:GLU:H	2.08	0.57
1:A:3:ALA:O	1:A:5:ALA:N	2.37	0.57
1:A:225:ILE:HD11	1:A:273:TRP:HZ3	1.70	0.57
1:A:113:LEU:HD13	1:A:119:LEU:HD11	1.86	0.56
1:A:249:ASP:OD2	4:A:605:HOH:O	2.18	0.56
1:A:333:LEU:C	1:A:335:ASP:H	2.07	0.56
1:A:13:ARG:HG3	1:A:53:LYS:HD3	1.85	0.56
1:A:202:GLY:HA2	4:A:514:HOH:O	2.05	0.56
1:A:22:ARG:HH11	1:A:44:LEU:HB3	1.70	0.56
1:A:178:HIS:CE1	1:A:199:ASN:HB2	2.40	0.56
1:A:302:VAL:HG13	1:A:303:GLU:N	2.20	0.55
1:A:35:GLY:HA3	1:A:53:LYS:O	2.06	0.55
1:A:2:ALA:O	1:A:4:ALA:N	2.35	0.55
1:A:9:PRO:C	1:A:11:MET:H	2.09	0.55
1:A:10:GLU:O	1:A:11:MET:HB2	2.06	0.55
1:A:133:ARG:HG3	1:A:320:GLU:HG2	1.89	0.55
1:A:322:ILE:HD11	4:A:709:HOH:O	2.05	0.55
1:A:17:PHE:CD2	1:A:36:MET:HE3	2.42	0.54
1:A:9:PRO:CB	1:A:15:GLN:OE1	2.55	0.54
1:A:125:CYS:HG	1:A:314:TYR:HE2	1.56	0.54
1:A:29:ILE:HD11	1:A:39:SER:N	2.23	0.54
1:A:86:ASP:HB3	1:A:102:VAL:HB	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:SER:O	1:A:222:ASN:HB2	2.06	0.54
1:A:9:PRO:HA	1:A:14:GLY:C	2.28	0.54
3:A:581:ATP:PB	3:A:581:ATP:H3'	2.48	0.53
1:A:328:LYS:HZ2	1:A:328:LYS:HB2	1.74	0.53
1:A:89:ARG:HD3	1:A:349:THR:OG1	2.08	0.53
1:A:336:LEU:HD12	1:A:341:LEU:HD21	1.89	0.53
1:A:13:ARG:HE	1:A:13:ARG:N	2.06	0.52
1:A:192:ARG:HD3	1:A:196:ILE:HG21	1.92	0.52
1:A:69:GLU:O	1:A:73:LEU:HB2	2.10	0.52
1:A:92:THR:HB	4:A:507:HOH:O	2.09	0.52
1:A:5:ALA:HB3	1:A:18:ASP:HA	1.90	0.52
1:A:74:LEU:HD13	1:A:84:ILE:HD12	1.91	0.52
1:A:202:GLY:HA2	1:A:207:ILE:HD13	1.92	0.52
1:A:59:HIS:ND1	1:A:62:TYR:CZ	2.78	0.51
1:A:13:ARG:HG3	1:A:53:LYS:CD	2.40	0.51
1:A:235:LEU:HD12	1:A:235:LEU:O	2.10	0.51
1:A:88:ILE:HB	1:A:100:TYR:HB2	1.91	0.51
1:A:340:LYS:HE3	1:A:344:LEU:HD13	1.93	0.51
1:A:185:TYR:CE2	1:A:187:ALA:CB	2.91	0.50
1:A:328:LYS:NZ	1:A:328:LYS:HB2	2.26	0.50
1:A:23:TYR:HE2	1:A:88:ILE:HD11	1.75	0.50
1:A:59:HIS:ND1	1:A:62:TYR:CE1	2.80	0.50
1:A:182:LEU:HA	1:A:231:TYR:HE2	1.76	0.50
1:A:204:THR:O	1:A:207:ILE:HB	2.12	0.50
1:A:330:ASP:O	1:A:331:MET:CB	2.58	0.50
1:A:329:PHE:CD1	1:A:329:PHE:O	2.65	0.49
1:A:333:LEU:C	1:A:335:ASP:N	2.66	0.49
1:A:43:ASN:HA	1:A:46:LYS:HZ3	1.75	0.49
1:A:9:PRO:CA	1:A:14:GLY:C	2.81	0.49
1:A:302:VAL:CG1	1:A:303:GLU:N	2.76	0.49
1:A:333:LEU:HD13	1:A:334:ASP:H	1.78	0.49
1:A:281:ASP:OD1	1:A:281:ASP:C	2.52	0.48
1:A:35:GLY:CA	1:A:53:LYS:O	2.60	0.48
1:A:279:ASN:OD1	1:A:279:ASN:N	2.43	0.48
1:A:122:ASP:HB3	4:A:508:HOH:O	2.13	0.48
1:A:2:ALA:C	1:A:4:ALA:N	2.65	0.47
1:A:182:LEU:HA	1:A:231:TYR:CE2	2.49	0.47
1:A:340:LYS:NZ	1:A:344:LEU:CD1	2.71	0.47
1:A:13:ARG:HB2	1:A:53:LYS:HZ3	1.80	0.47
1:A:110:LEU:HB3	1:A:150:PRO:HB3	1.97	0.47
1:A:216:LEU:HD12	1:A:291:MET:HE1	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:LYS:NZ	1:A:328:LYS:HB3	2.30	0.47
1:A:67:LEU:O	1:A:71:LYS:HG3	2.15	0.47
1:A:67:LEU:HD22	1:A:71:LYS:HG3	1.97	0.47
1:A:65:ARG:HG3	1:A:167:GLY:CA	2.46	0.46
1:A:110:LEU:HD12	1:A:153:LEU:HD12	1.96	0.46
1:A:158:THR:O	1:A:159:CYS:HB2	2.15	0.46
1:A:216:LEU:HD12	1:A:291:MET:CE	2.46	0.46
1:A:328:LYS:HB3	1:A:328:LYS:HZ3	1.81	0.46
3:A:581:ATP:H3'	3:A:581:ATP:O2B	2.15	0.46
1:A:29:ILE:HG22	1:A:29:ILE:O	2.15	0.46
1:A:65:ARG:HE	1:A:65:ARG:HB2	1.56	0.46
1:A:29:ILE:CD1	1:A:39:SER:N	2.78	0.46
1:A:106:MET:O	3:A:581:ATP:C2	2.69	0.46
1:A:17:PHE:CD2	1:A:36:MET:HE2	2.47	0.46
1:A:22:ARG:HG2	1:A:23:TYR:CE2	2.51	0.46
1:A:239:LEU:O	1:A:243:GLY:N	2.49	0.46
1:A:230:HIS:NE2	1:A:233:ASP:HB2	2.31	0.45
1:A:57:PHE:CE1	1:A:342:LYS:HG3	2.52	0.45
1:A:104:ASP:O	3:A:581:ATP:N6	2.46	0.45
1:A:14:GLY:HA2	1:A:36:MET:HE1	1.98	0.45
1:A:202:GLY:HA2	1:A:207:ILE:CD1	2.47	0.45
1:A:250:LEU:HD12	1:A:250:LEU:HA	1.80	0.45
1:A:327:PHE:CD1	1:A:328:LYS:N	2.85	0.45
1:A:69:GLU:HG3	1:A:73:LEU:HD22	1.98	0.45
1:A:272:PRO:HG2	1:A:275:ARG:HG3	1.99	0.45
1:A:336:LEU:HD22	1:A:336:LEU:HA	1.65	0.44
1:A:43:ASN:HA	1:A:46:LYS:NZ	2.33	0.44
1:A:336:LEU:HD12	1:A:341:LEU:HD23	1.96	0.44
1:A:141:ALA:O	1:A:142:ASN:CB	2.65	0.44
1:A:106:MET:O	3:A:581:ATP:H2	2.01	0.44
1:A:12:VAL:HG13	1:A:33:ALA:HA	1.99	0.44
1:A:250:LEU:HD11	1:A:262:LEU:HD23	2.00	0.44
1:A:230:HIS:CD2	1:A:233:ASP:HB2	2.53	0.43
1:A:89:ARG:NH1	1:A:346:PHE:O	2.52	0.43
1:A:337:PRO:C	1:A:339:GLU:N	2.71	0.43
1:A:262:LEU:HD12	1:A:262:LEU:HA	1.84	0.43
1:A:59:HIS:HB3	1:A:62:TYR:CD2	2.53	0.43
1:A:281:ASP:OD1	1:A:283:LYS:N	2.40	0.43
1:A:273:TRP:HB3	1:A:285:LEU:HD22	2.01	0.43
1:A:189:ARG:O	1:A:192:ARG:HB2	2.19	0.42
1:A:333:LEU:HD13	1:A:334:ASP:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:GLU:O	1:A:222:ASN:N	2.49	0.42
1:A:108:THR:OG1	1:A:109:ASP:N	2.51	0.42
1:A:313:GLN:HB3	4:A:533:HOH:O	2.20	0.42
1:A:187:ALA:HB1	4:A:541:HOH:O	2.19	0.42
1:A:308:HIS:CG	1:A:309:PRO:HD2	2.55	0.42
1:A:115:LYS:HB2	1:A:115:LYS:HE2	1.73	0.42
1:A:344:LEU:HD12	1:A:344:LEU:HA	1.79	0.42
1:A:22:ARG:HG2	1:A:23:TYR:CE1	2.55	0.41
1:A:68:ARG:HD3	1:A:169:ALA:O	2.20	0.41
1:A:78:HIS:CB	1:A:81:ILE:HD12	2.49	0.41
1:A:114:LEU:HD13	1:A:218:GLU:HG2	2.02	0.41
1:A:333:LEU:HD22	1:A:333:LEU:HA	1.92	0.41
1:A:213:GLY:HA2	1:A:291:MET:HE3	2.02	0.41
1:A:67:LEU:O	1:A:67:LEU:HD22	2.21	0.41
1:A:54:ILE:HG22	1:A:56:PRO:CD	2.51	0.40
1:A:87:ILE:HG21	1:A:349:THR:HG22	2.03	0.40
1:A:277:PHE:HA	1:A:278:PRO:HD2	1.80	0.40
1:A:70:ILE:HG23	1:A:74:LEU:HD22	2.03	0.40
1:A:87:ILE:HD12	1:A:352:PHE:CE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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	355/364 (98%)	315 (89%)	27 (8%)	13 (4%)	3 11

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	331	MET
1	A	332	GLU

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Mol	Chain	Res	Type
1	A	357	ARG
1	A	4	ALA
1	A	8	GLY
1	A	11	MET
1	A	25	ASN
1	A	12	VAL
1	A	15	GLN
1	A	24	THR
1	A	31	GLU
1	A	9	PRO
1	A	354	PRO

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	317/324 (98%)	264 (83%)	53 (17%)	2 6

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	MET
1	A	13	ARG
1	A	15	GLN
1	A	22	ARG
1	A	24	THR
1	A	31	GLU
1	A	39	SER
1	A	46	LYS
1	A	55	SER
1	A	59	HIS
1	A	60	GLN
1	A	67	LEU
1	A	73	LEU
1	A	74	LEU
1	A	75	ARG

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Mol	Chain	Res	Type
1	A	82	ILE
1	A	88	ILE
1	A	89	ARG
1	A	93	ILE
1	A	98	ASP
1	A	107	GLU
1	A	110	LEU
1	A	113	LEU
1	A	115	LYS
1	A	118	HIS
1	A	142	ASN
1	A	155	LEU
1	A	158	THR
1	A	184	GLU
1	A	196	ILE
1	A	199	ASN
1	A	207	ILE
1	A	216	LEU
1	A	229	LYS
1	A	237	HIS
1	A	242	LEU
1	A	247	GLN
1	A	250	LEU
1	A	255	ASN
1	A	262	LEU
1	A	265	LEU
1	A	279	ASN
1	A	303	GLU
1	A	322	ILE
1	A	329	PHE
1	A	331	MET
1	A	333	LEU
1	A	336	LEU
1	A	340	LYS
1	A	344	LEU
1	A	347	GLU
1	A	353	GLN
1	A	357	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	43	ASN
1	A	45	ASN
1	A	78	HIS
1	A	121	ASN
1	A	152	ASN
1	A	178	HIS
1	A	199	ASN
1	A	251	ASN
1	A	255	ASN
1	A	260	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\#\|Z\| > 2$
3	ATP	A	581	2	26,33,33	7.97	11 (42%)	31,52,52	2.92	13 (41%)

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
3	ATP	A	581	2	-	6/18/38/38	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	581	ATP	C8-N7	33.29	1.94	1.34
3	A	581	ATP	O4'-C1'	17.76	1.65	1.41
3	A	581	ATP	O4'-C4'	7.44	1.61	1.45
3	A	581	ATP	PA-O5'	6.99	1.87	1.59
3	A	581	ATP	C2-N1	6.20	1.45	1.33
3	A	581	ATP	C5'-C4'	4.94	1.67	1.51
3	A	581	ATP	C3'-C4'	4.13	1.63	1.53
3	A	581	ATP	O5'-C5'	3.22	1.57	1.44
3	A	581	ATP	C2'-C3'	-2.71	1.45	1.53
3	A	581	ATP	C2'-C1'	2.56	1.57	1.53
3	A	581	ATP	C6-N1	2.41	1.47	1.37

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	581	ATP	C5-C6-N1	-7.18	104.07	120.35
3	A	581	ATP	C5-C6-N6	6.40	130.07	120.35
3	A	581	ATP	C1'-N9-C4	-6.16	115.83	126.64
3	A	581	ATP	O5'-C5'-C4'	4.47	124.37	108.99
3	A	581	ATP	C2-N1-C6	4.10	125.77	118.75
3	A	581	ATP	C5'-C4'-C3'	3.76	129.27	115.18
3	A	581	ATP	C3'-C2'-C1'	3.71	106.56	100.98
3	A	581	ATP	N3-C2-N1	-3.58	123.08	128.68
3	A	581	ATP	N6-C6-N1	3.48	125.79	118.57
3	A	581	ATP	O3'-C3'-C4'	-3.03	102.30	111.05
3	A	581	ATP	O2'-C2'-C3'	2.37	119.49	111.82
3	A	581	ATP	C4-C5-N7	2.36	111.86	109.40
3	A	581	ATP	O3G-PG-O2G	2.28	116.36	107.64

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	581	ATP	C5'-O5'-PA-O3A
3	A	581	ATP	C4'-C5'-O5'-PA
3	A	581	ATP	C5'-O5'-PA-O1A
3	A	581	ATP	C5'-O5'-PA-O2A
3	A	581	ATP	C3'-C4'-C5'-O5'
3	A	581	ATP	PB-O3B-PG-O3G

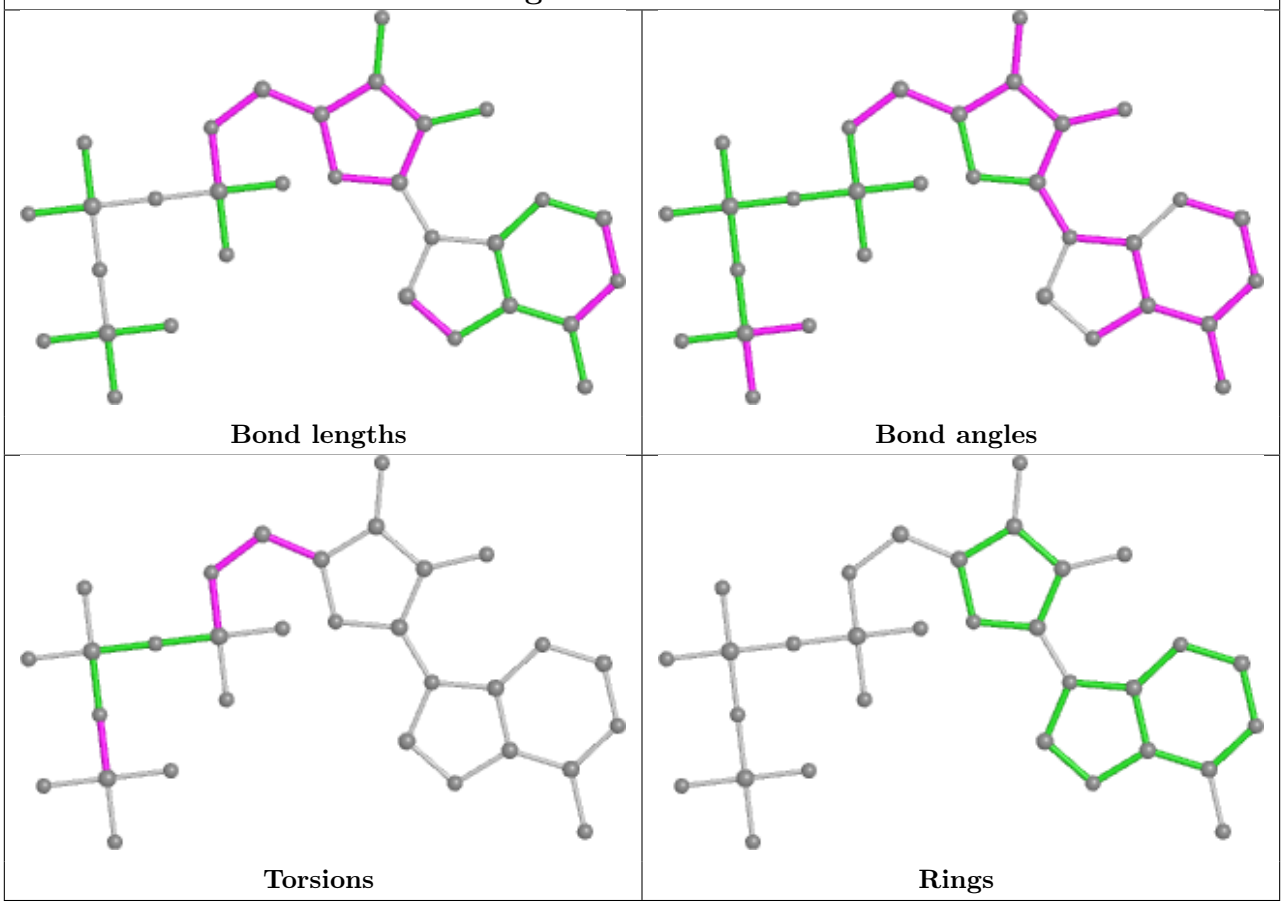
There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	581	ATP	7	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 then 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.

Ligand ATP A 581



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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

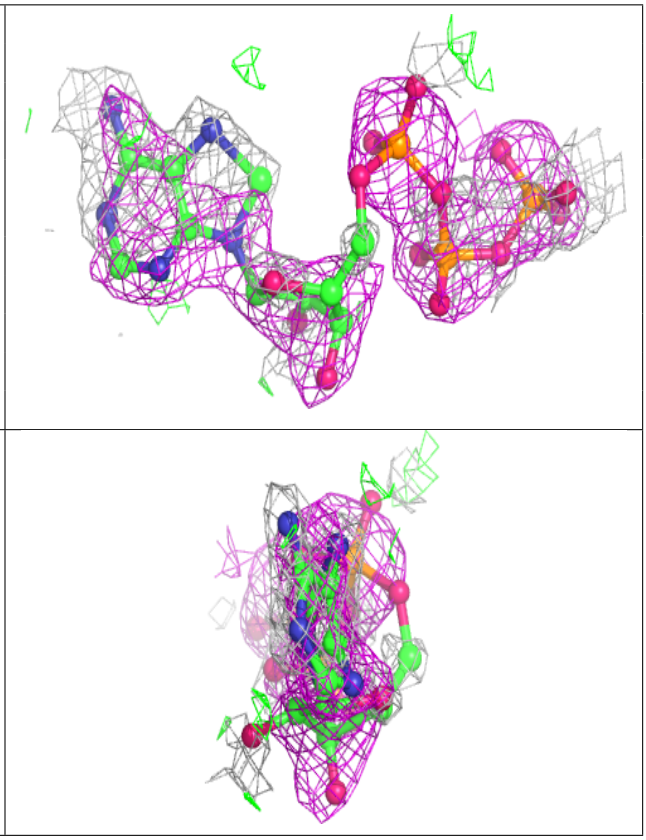
6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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 A 581:

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.