



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:39 PM UTC

PDB ID : 4YC4 / pdb_00004yc4
Title : Crystal structure of phosphatidyl inositol 4-kinase II alpha in complex with nucleotide analog
Authors : Klima, M.; Boura, E.
Deposited on : 2015-02-19
Resolution : 2.58 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

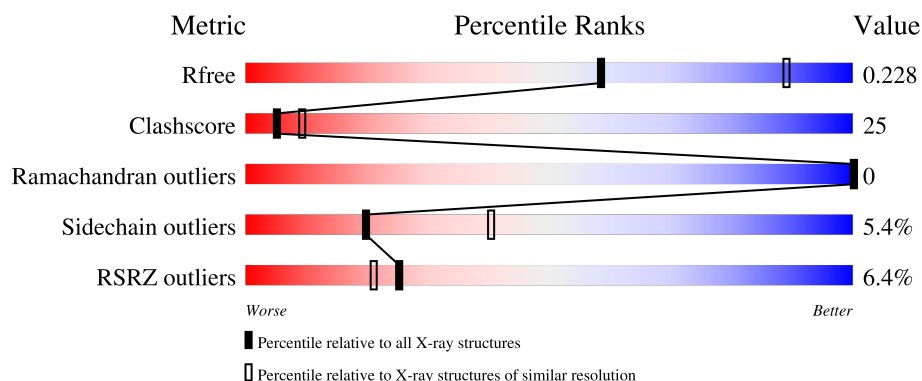
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	556	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3943 atoms, of which 12 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.

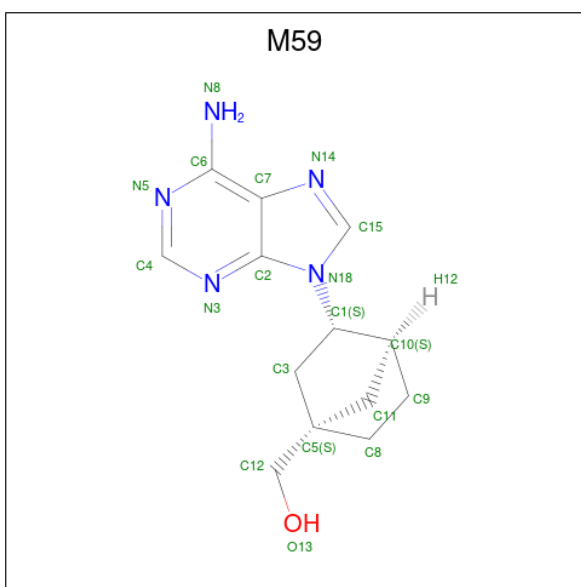
- Molecule 1 is a protein called Phosphatidylinositol 4-kinase type 2-alpha, Lysozyme, Phosphatidylinositol 4-kinase type 2-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	0	0
			3882	2490	674	710	8			

There are 9 discrepancies between the modelled and reference sequences:

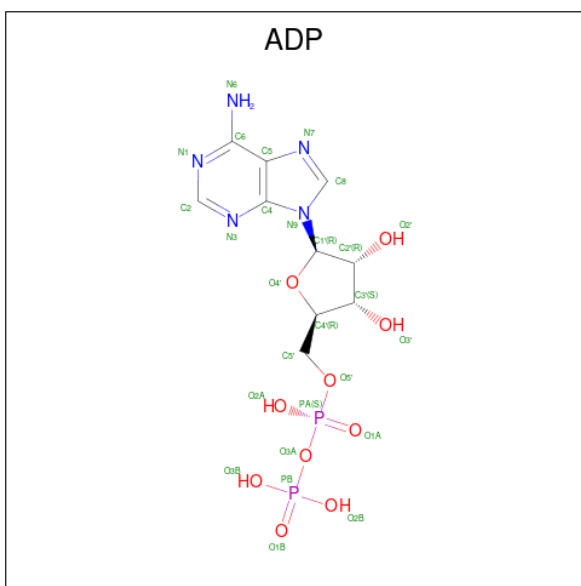
Chain	Residue	Modelled	Actual	Comment	Reference
A	72	GLY	-	expression tag	UNP Q9BTU6
A	73	ALA	-	expression tag	UNP Q9BTU6
A	74	MET	-	expression tag	UNP Q9BTU6
A	75	GLY	-	expression tag	UNP Q9BTU6
A	1001	GLY	-	linker	UNP Q9BTU6
A	1002	THR	-	linker	UNP Q9BTU6
A	1003	GLY	-	linker	UNP Q9BTU6
A	1056	THR	CYS	conflict	UNP D9IEF7
A	1099	ALA	CYS	conflict	UNP D9IEF7

- Molecule 2 is [(1S,3S,4S)-3-(6-amino-9H-purin-9-yl)bicyclo[2.2.1]hept-1-yl]methanol (CCD ID: M59) (formula: C₁₃H₁₇N₅O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			19	13	5	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	H	N	O	P	0	0
			39	10	12	5	10	2		

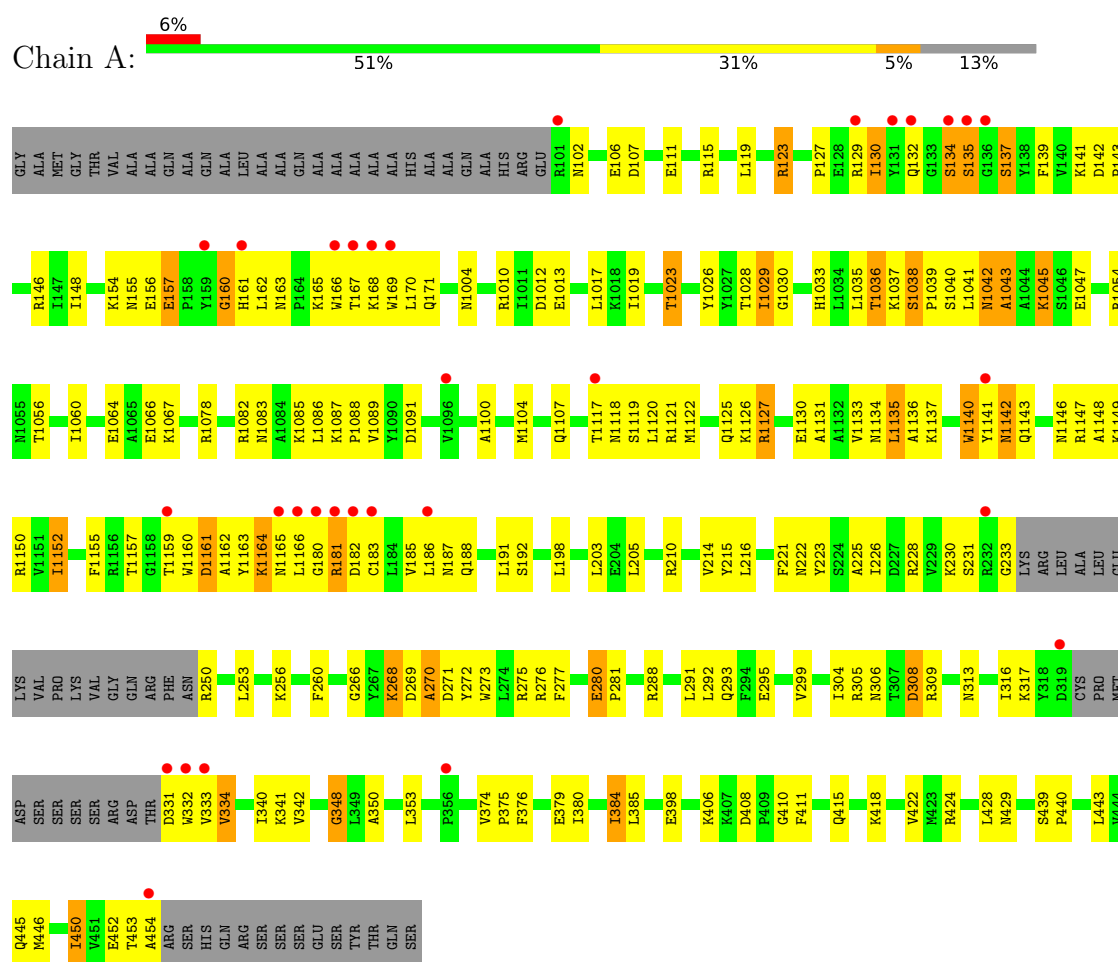
- Molecule 4 is water.

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

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: Phosphatidylinositol 4-kinase type 2-alpha, Lysozyme, Phosphatidylinositol 4-kinase type 2-alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	79.41Å 103.77Å 78.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.07 – 2.58 49.07 – 2.58	Depositor EDS
% Data completeness (in resolution range)	95.8 (49.07-2.58) 91.1 (49.07-2.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.214 , 0.256 (Not available) , 0.228	Depositor DCC
R_{free} test set	1007 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.907	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3943	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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: M59, ADP

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	0.72	0/3971	1.13	25/5387 (0.5%)

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

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

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	GLY	N-CA-C	8.46	123.97	114.67
1	A	135	SER	N-CA-C	8.13	121.99	112.72
1	A	374	VAL	N-CA-C	7.63	115.20	108.63
1	A	1140	TRP	N-CA-C	-7.06	101.73	110.41
1	A	1038	SER	CA-C-N	7.00	126.63	119.56
1	A	1038	SER	C-N-CA	7.00	126.63	119.56
1	A	137	SER	N-CA-C	6.80	120.32	109.24
1	A	1166	LEU	N-CA-C	6.72	119.81	107.99
1	A	132	GLN	N-CA-C	6.21	119.95	112.38
1	A	1163	TYR	N-CA-C	6.01	120.74	113.17
1	A	1164	LYS	N-CA-C	5.93	120.79	113.50
1	A	308	ASP	N-CA-C	5.92	118.45	110.88
1	A	270	ALA	N-CA-C	-5.83	104.11	111.11
1	A	1161	ASP	N-CA-C	5.72	118.25	111.33
1	A	134	SER	N-CA-C	-5.56	102.79	110.35
1	A	1043	ALA	N-CA-C	-5.50	104.98	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ARG	N-CA-C	-5.48	99.13	110.80
1	A	1045	LYS	N-CA-C	-5.33	105.55	111.36
1	A	186	LEU	N-CA-C	5.29	116.85	111.14
1	A	384	ILE	N-CA-C	5.18	120.11	109.34
1	A	280	GLU	CA-C-N	5.11	125.02	119.76
1	A	280	GLU	C-N-CA	5.11	125.02	119.76
1	A	253	LEU	CA-C-N	-5.02	116.05	119.66
1	A	253	LEU	C-N-CA	-5.02	116.05	119.66
1	A	160	GLY	N-CA-C	5.01	118.36	111.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	134	SER	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3882	0	3847	185	1
2	A	19	0	17	3	0
3	A	27	12	12	7	0
4	A	3	0	0	0	0
All	All	3931	12	3876	192	1

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 25.

All (192) 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:332:TRP:NE1	1:A:334:VAL:HG13	1.53	1.24
1:A:332:TRP:HE1	1:A:334:VAL:CG1	1.60	1.15
1:A:453:THR:HG22	1:A:454:ALA:H	1.09	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:VAL:HG11	1:A:384:ILE:HD13	1.40	1.02
1:A:123:ARG:HH11	1:A:123:ARG:HG3	1.20	1.01
1:A:332:TRP:CD1	1:A:334:VAL:H	1.78	1.01
1:A:398:GLU:OE1	1:A:424:ARG:NH2	1.94	0.99
1:A:453:THR:HG22	1:A:454:ALA:N	1.76	0.99
1:A:453:THR:CG2	1:A:454:ALA:H	1.74	0.98
1:A:272:TYR:CZ	1:A:276:ARG:NH2	2.33	0.96
1:A:332:TRP:HE1	1:A:334:VAL:HG13	0.75	0.91
1:A:1157:THR:OG1	1:A:1159:THR:HG22	1.71	0.89
1:A:332:TRP:HD1	1:A:334:VAL:H	0.93	0.89
1:A:332:TRP:HD1	1:A:334:VAL:N	1.72	0.86
1:A:123:ARG:HH11	1:A:123:ARG:CG	1.90	0.84
1:A:332:TRP:CD1	1:A:334:VAL:N	2.46	0.83
1:A:429:ASN:CG	1:A:446:MET:HE2	2.04	0.83
1:A:1023:THR:HG21	1:A:1143:GLN:HG3	1.62	0.81
1:A:272:TYR:CE1	1:A:276:ARG:NH2	2.51	0.79
1:A:1133:VAL:O	1:A:1136:ALA:HB3	1.85	0.77
1:A:332:TRP:CD1	1:A:333:VAL:N	2.53	0.76
1:A:1054:ARG:HH12	1:A:1056:THR:HA	1.50	0.76
1:A:166:TRP:CH2	1:A:1164:LYS:CB	2.70	0.75
1:A:1122:MET:HB2	1:A:1131:ALA:HB2	1.70	0.73
1:A:1054:ARG:NH1	1:A:1056:THR:HA	2.03	0.73
1:A:331:ASP:OD1	1:A:331:ASP:O	2.08	0.72
1:A:1140:TRP:O	1:A:1141:TYR:HB3	1.89	0.71
1:A:1141:TYR:OH	1:A:1149:LYS:HE2	1.90	0.71
1:A:1165:ASN:O	1:A:180:GLY:HA2	1.91	0.71
1:A:107:ASP:OD2	1:A:210:ARG:NH1	2.25	0.70
1:A:1119:SER:HB3	1:A:1135:LEU:HD11	1.74	0.70
1:A:1012:ASP:OD1	1:A:1150:ARG:NH1	2.24	0.70
1:A:299:VAL:HG11	1:A:384:ILE:CD1	2.20	0.69
2:A:1201:M59:C15	2:A:1201:M59:H13	2.23	0.69
1:A:1118:ASN:HA	1:A:1121:ARG:HH12	1.58	0.68
3:A:1202:ADP:H3'	3:A:1202:ADP:O2A	1.93	0.68
1:A:155:ASN:HD21	1:A:256:LYS:HE2	1.58	0.68
1:A:1013:GLU:OE2	1:A:1147:ARG:NH1	2.23	0.68
1:A:192:SER:HB3	1:A:348:GLY:HA2	1.76	0.68
1:A:205:LEU:HD21	1:A:384:ILE:HD11	1.76	0.67
1:A:1148:ALA:O	1:A:1152:ILE:HG13	1.94	0.66
1:A:119:LEU:HD11	1:A:123:ARG:NH1	2.10	0.66
2:A:1201:M59:H13	2:A:1201:M59:H4	1.79	0.65
1:A:268:LYS:HD3	1:A:272:TYR:CE1	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:LEU:C	1:A:428:LEU:HD23	2.23	0.64
1:A:445:GLN:HG2	3:A:1202:ADP:H1'	1.78	0.64
1:A:308:ASP:OD1	1:A:308:ASP:O	2.14	0.64
1:A:1041:LEU:CD2	1:A:1045:LYS:HE3	2.28	0.64
1:A:353:LEU:HD21	1:A:415:GLN:HG2	1.80	0.64
1:A:1135:LEU:CD1	1:A:1135:LEU:N	2.61	0.63
1:A:1023:THR:CG2	1:A:1143:GLN:HG3	2.28	0.63
3:A:1202:ADP:H3'	3:A:1202:ADP:PA	2.38	0.63
1:A:453:THR:CG2	1:A:454:ALA:N	2.42	0.63
1:A:1029:ILE:CD1	1:A:1060:ILE:HG12	2.29	0.63
1:A:129:ARG:HD2	1:A:135:SER:O	1.98	0.62
1:A:1119:SER:CB	1:A:1135:LEU:HD11	2.29	0.62
1:A:226:ILE:HG22	1:A:230:LYS:HE2	1.82	0.61
1:A:1086:LEU:HD23	1:A:1120:LEU:HD21	1.81	0.61
1:A:1041:LEU:O	1:A:1045:LYS:HG3	2.01	0.61
1:A:1140:TRP:O	1:A:1142:ASN:N	2.33	0.60
1:A:331:ASP:OD1	1:A:331:ASP:C	2.44	0.60
1:A:1029:ILE:HD13	1:A:1060:ILE:CD1	2.32	0.60
1:A:1028:THR:HG22	1:A:1029:ILE:N	2.17	0.59
1:A:228:ARG:O	1:A:231:SER:OG	2.19	0.59
1:A:291:LEU:O	1:A:295:GLU:HG3	2.02	0.59
1:A:123:ARG:CG	1:A:123:ARG:NH1	2.57	0.58
1:A:1100:ALA:HB3	1:A:1155:PHE:CE1	2.38	0.58
1:A:380:ILE:HD11	1:A:384:ILE:HD12	1.85	0.58
1:A:268:LYS:HD3	1:A:272:TYR:CD1	2.39	0.57
1:A:272:TYR:OH	1:A:276:ARG:NH2	2.37	0.57
1:A:277:PHE:O	1:A:281:PRO:HA	2.05	0.57
1:A:266:GLY:O	1:A:317:LYS:HG3	2.05	0.57
1:A:1087:LYS:HG2	1:A:1091:ASP:OD2	2.05	0.56
1:A:1035:LEU:O	1:A:1047:GLU:HG2	2.05	0.56
1:A:308:ASP:O	1:A:308:ASP:CG	2.48	0.56
1:A:308:ASP:O	1:A:313:ASN:ND2	2.39	0.56
1:A:130:ILE:HG13	1:A:137:SER:O	2.06	0.56
1:A:1119:SER:OG	1:A:1135:LEU:CD1	2.54	0.56
1:A:1040:SER:OG	1:A:1043:ALA:HB2	2.06	0.56
1:A:1085:LYS:HZ3	1:A:1117:THR:HG22	1.71	0.55
1:A:1041:LEU:HD21	1:A:1045:LYS:CE	2.36	0.55
1:A:1118:ASN:HA	1:A:1121:ARG:NH1	2.21	0.55
1:A:1085:LYS:NZ	1:A:1117:THR:HG22	2.22	0.54
1:A:333:VAL:HG12	1:A:333:VAL:O	2.07	0.54
1:A:1119:SER:OG	1:A:1134:ASN:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1029:ILE:HD11	1:A:1060:ILE:HG12	1.90	0.53
1:A:1041:LEU:O	1:A:1041:LEU:HG	2.08	0.53
1:A:163:ASN:OD1	1:A:165:LYS:N	2.41	0.53
1:A:408:ASP:OD1	1:A:410:GLY:N	2.40	0.53
1:A:440:PRO:O	1:A:443:LEU:HB3	2.10	0.52
1:A:305:ARG:O	1:A:305:ARG:HG3	2.09	0.52
1:A:1041:LEU:HD21	1:A:1045:LYS:HE3	1.91	0.52
1:A:119:LEU:CD1	1:A:123:ARG:NH1	2.73	0.52
1:A:1135:LEU:N	1:A:1135:LEU:HD12	2.24	0.52
1:A:304:ILE:HD12	1:A:350:ALA:CB	2.39	0.52
1:A:154:LYS:O	1:A:157:GLU:HB2	2.10	0.52
1:A:1159:THR:HG23	1:A:1161:ASP:H	1.75	0.52
1:A:1140:TRP:C	1:A:1142:ASN:H	2.15	0.52
1:A:1029:ILE:HD13	1:A:1060:ILE:HG12	1.91	0.51
1:A:162:LEU:O	1:A:222:ASN:ND2	2.41	0.51
1:A:155:ASN:ND2	1:A:256:LYS:HE2	2.25	0.51
1:A:271:ASP:OD2	1:A:275:ARG:NH2	2.42	0.51
1:A:332:TRP:CD1	1:A:333:VAL:H	2.29	0.51
1:A:269:ASP:O	1:A:270:ALA:C	2.53	0.51
1:A:445:GLN:CG	3:A:1202:ADP:HI'	2.42	0.50
1:A:1141:TYR:HA	1:A:1148:ALA:CB	2.41	0.50
1:A:215:TYR:O	1:A:216:LEU:HD23	2.11	0.50
1:A:141:LYS:HA	1:A:146:ARG:O	2.12	0.49
1:A:1140:TRP:C	1:A:1142:ASN:N	2.70	0.49
1:A:1087:LYS:N	1:A:1088:PRO:CD	2.76	0.49
1:A:406:LYS:HA	1:A:411:PHE:CD2	2.47	0.49
1:A:1121:ARG:NH1	1:A:1121:ARG:HB2	2.28	0.49
1:A:182:ASP:O	1:A:182:ASP:OD1	2.31	0.49
1:A:1086:LEU:HD23	1:A:1120:LEU:CD2	2.44	0.48
3:A:1202:ADP:PA	3:A:1202:ADP:C3'	3.01	0.48
1:A:181:ARG:C	1:A:183:CYS:H	2.21	0.48
1:A:1118:ASN:O	1:A:1121:ARG:HB3	2.13	0.48
1:A:1127:ARG:HD3	1:A:1130:GLU:OE2	2.13	0.48
1:A:293:GLN:OE1	1:A:341:LYS:HD2	2.13	0.48
1:A:406:LYS:HA	1:A:411:PHE:CG	2.48	0.48
1:A:156:GLU:OE2	1:A:256:LYS:NZ	2.47	0.48
1:A:1083:ASN:OD1	1:A:1085:LYS:N	2.47	0.48
1:A:168:LYS:HG3	1:A:169:TRP:N	2.28	0.47
1:A:1083:ASN:OD1	1:A:1083:ASN:C	2.57	0.47
1:A:1104:MET:HE3	1:A:1140:TRP:CZ3	2.49	0.47
1:A:233:GLY:C	1:A:250:ARG:NH1	2.72	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1010:ARG:HH21	1:A:1066:GLU:CD	2.22	0.47
1:A:119:LEU:HD23	1:A:143:PRO:HG2	1.96	0.47
1:A:271:ASP:OD1	1:A:275:ARG:NH2	2.48	0.47
1:A:428:LEU:HD23	1:A:428:LEU:O	2.15	0.47
1:A:1085:LYS:NZ	1:A:1117:THR:CG2	2.78	0.47
3:A:1202:ADP:O2A	3:A:1202:ADP:C3'	2.62	0.47
1:A:1029:ILE:HG12	1:A:1030:GLY:H	1.80	0.47
1:A:198:LEU:C	1:A:198:LEU:HD23	2.40	0.46
1:A:123:ARG:HG3	1:A:123:ARG:NH1	2.02	0.46
1:A:1029:ILE:HD13	1:A:1060:ILE:HD13	1.97	0.46
1:A:106:GLU:HG2	1:A:107:ASP:N	2.30	0.45
1:A:191:LEU:HD11	1:A:408:ASP:HB2	1.98	0.45
2:A:1201:M59:C15	2:A:1201:M59:C11	2.93	0.45
1:A:1159:THR:HG21	1:A:1161:ASP:OD2	2.16	0.45
1:A:288:ARG:CZ	1:A:292:LEU:HD21	2.45	0.45
1:A:1004:ASN:HB2	1:A:1160:TRP:CZ2	2.51	0.45
1:A:1029:ILE:HD13	1:A:1060:ILE:CG1	2.47	0.45
1:A:418:LYS:O	1:A:422:VAL:HG23	2.16	0.45
1:A:1125:GLN:O	1:A:1126:LYS:HB2	2.17	0.45
1:A:280:GLU:O	1:A:280:GLU:HG3	2.17	0.45
1:A:332:TRP:CG	1:A:333:VAL:N	2.84	0.45
1:A:1127:ARG:HD3	1:A:1130:GLU:CD	2.42	0.44
1:A:157:GLU:O	1:A:160:GLY:N	2.46	0.44
1:A:203:LEU:HB2	1:A:205:LEU:HG	1.99	0.44
1:A:225:ALA:O	1:A:226:ILE:C	2.60	0.44
1:A:1078:ARG:CZ	1:A:1082:ARG:HD2	2.47	0.44
1:A:1089:VAL:HG21	1:A:1120:LEU:HB3	2.00	0.44
1:A:271:ASP:CG	1:A:275:ARG:NH2	2.76	0.44
1:A:130:ILE:HG13	1:A:137:SER:C	2.43	0.44
1:A:332:TRP:CD1	1:A:334:VAL:HG13	2.42	0.44
1:A:214:VAL:HB	1:A:260:PHE:HB2	2.00	0.44
1:A:1150:ARG:HG3	1:A:1162:ALA:HB1	2.00	0.44
1:A:127:PRO:HA	1:A:139:PHE:O	2.18	0.43
1:A:1026:TYR:CZ	1:A:1037:LYS:HD3	2.54	0.43
1:A:1036:THR:OG1	1:A:1038:SER:HB3	2.18	0.43
1:A:1042:ASN:OD1	1:A:1042:ASN:N	2.51	0.43
1:A:332:TRP:CD1	1:A:332:TRP:C	2.96	0.43
1:A:1121:ARG:NH1	1:A:1121:ARG:CB	2.82	0.43
1:A:161:HIS:O	1:A:161:HIS:CD2	2.72	0.43
1:A:306:ASN:ND2	1:A:309:ARG:HB2	2.34	0.43
1:A:385:LEU:HD23	1:A:385:LEU:HA	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1064:GLU:O	1:A:1067:LYS:HG2	2.19	0.42
1:A:445:GLN:HG2	3:A:1202:ADP:C1'	2.45	0.42
1:A:1135:LEU:H	1:A:1135:LEU:HD13	1.84	0.42
1:A:142:ASP:HB3	1:A:148:ILE:HD13	2.01	0.42
1:A:1107:GLN:OE1	1:A:1140:TRP:NE1	2.52	0.42
1:A:273:TRP:CD1	1:A:316:ILE:HB	2.55	0.42
1:A:276:ARG:CZ	1:A:276:ARG:HB2	2.50	0.42
1:A:1120:LEU:HD12	1:A:1120:LEU:N	2.35	0.42
1:A:1120:LEU:HD12	1:A:1120:LEU:H	1.84	0.41
1:A:450:ILE:HD12	1:A:452:GLU:HB2	2.02	0.41
1:A:1038:SER:HA	1:A:1039:PRO:HD3	1.74	0.41
1:A:111:GLU:O	1:A:115:ARG:HG2	2.20	0.41
1:A:1104:MET:HE3	1:A:1140:TRP:CH2	2.56	0.41
1:A:102:ASN:O	1:A:115:ARG:NH1	2.48	0.41
1:A:185:VAL:HG11	1:A:188:GLN:CD	2.45	0.41
1:A:1028:THR:CG2	1:A:1029:ILE:N	2.83	0.41
1:A:1028:THR:HG23	1:A:1033:HIS:O	2.20	0.41
1:A:1033:HIS:CD2	1:A:1033:HIS:C	2.98	0.41
1:A:221:PHE:HB2	1:A:223:TYR:CE2	2.55	0.41
1:A:375:PRO:O	1:A:376:PHE:C	2.64	0.41
1:A:428:LEU:C	1:A:428:LEU:CD2	2.92	0.41
1:A:1019:ILE:CD1	1:A:1045:LYS:HG2	2.51	0.41
1:A:1017:LEU:HD23	1:A:1017:LEU:HA	1.76	0.40
1:A:187:ASN:O	1:A:187:ASN:CG	2.64	0.40
1:A:185:VAL:HG11	1:A:188:GLN:OE1	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1078:ARG:NH2	1:A:379:GLU:OE2[1_556]	2.08	0.12

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	480/556 (86%)	463 (96%)	17 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	409/473 (86%)	387 (95%)	22 (5%)	20	40

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	ARG
1	A	130	ILE
1	A	157	GLU
1	A	167	THR
1	A	170	LEU
1	A	171	GLN
1	A	1023	THR
1	A	1029	ILE
1	A	1036	THR
1	A	1042	ASN
1	A	1127	ARG
1	A	1135	LEU
1	A	1137	LYS
1	A	1142	ASN
1	A	1146	ASN
1	A	1152	ILE
1	A	268	LYS
1	A	334	VAL
1	A	340	ILE
1	A	342	VAL
1	A	439	SER

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Mol	Chain	Res	Type
1	A	450	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	161	HIS
1	A	1070	ASN
1	A	1124	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](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	M59	A	1201	-	22,22,22	1.89	4 (18%)	26,34,34	2.51	11 (42%)
3	ADP	A	1202	-	28,29,29	1.31	4 (14%)	43,45,45	2.06	10 (23%)

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	M59	A	1201	-	-	3/7/27/27	0/5/4/4
3	ADP	A	1202	-	-	5/16/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	M59	C6-N8	6.29	1.50	1.34
3	A	1202	ADP	C5-C4	4.10	1.46	1.39
2	A	1201	M59	C11-C5	-3.25	1.51	1.54
2	A	1201	M59	C1-N18	2.74	1.49	1.47
2	A	1201	M59	C15-N14	2.30	1.36	1.31
3	A	1202	ADP	C8-N7	2.22	1.35	1.31
3	A	1202	ADP	C5-N7	-2.04	1.35	1.39
3	A	1202	ADP	C5-C6	2.04	1.46	1.41

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1202	ADP	C5-C4-N3	-6.06	118.37	126.72
2	A	1201	M59	C7-C2-N3	-5.75	118.79	126.72
3	A	1202	ADP	N3-C4-N9	5.52	136.56	127.17
2	A	1201	M59	N5-C4-N3	-4.37	121.97	128.58
3	A	1202	ADP	C2-N3-C4	4.24	122.18	111.83
2	A	1201	M59	C3-C5-C12	4.21	116.60	110.71
3	A	1202	ADP	N3-C2-N1	-4.10	122.38	128.58
3	A	1202	ADP	C4-N9-C8	3.69	109.61	105.74
2	A	1201	M59	C4-N3-C2	3.49	120.36	111.83
2	A	1201	M59	C3-C1-N18	-3.44	108.95	113.30
2	A	1201	M59	C2-C7-N14	-3.38	106.72	110.58
2	A	1201	M59	N3-C2-N18	3.19	132.59	127.17
2	A	1201	M59	C11-C5-C12	3.15	115.12	110.71
3	A	1202	ADP	C2'-C3'-C4'	2.93	108.27	102.61
2	A	1201	M59	N18-C15-N14	-2.84	109.90	113.94
2	A	1201	M59	C7-N14-C15	2.66	107.62	103.45
3	A	1202	ADP	N9-C8-N7	-2.59	110.26	113.94
2	A	1201	M59	C7-C2-N18	2.56	108.60	105.81
3	A	1202	ADP	C5-C6-N6	-2.31	117.58	123.29
3	A	1202	ADP	C5-N7-C8	2.30	107.07	103.45
3	A	1202	ADP	C4-C5-N7	-2.26	107.99	110.58

There are no chirality outliers.

All (8) torsion outliers are listed below:

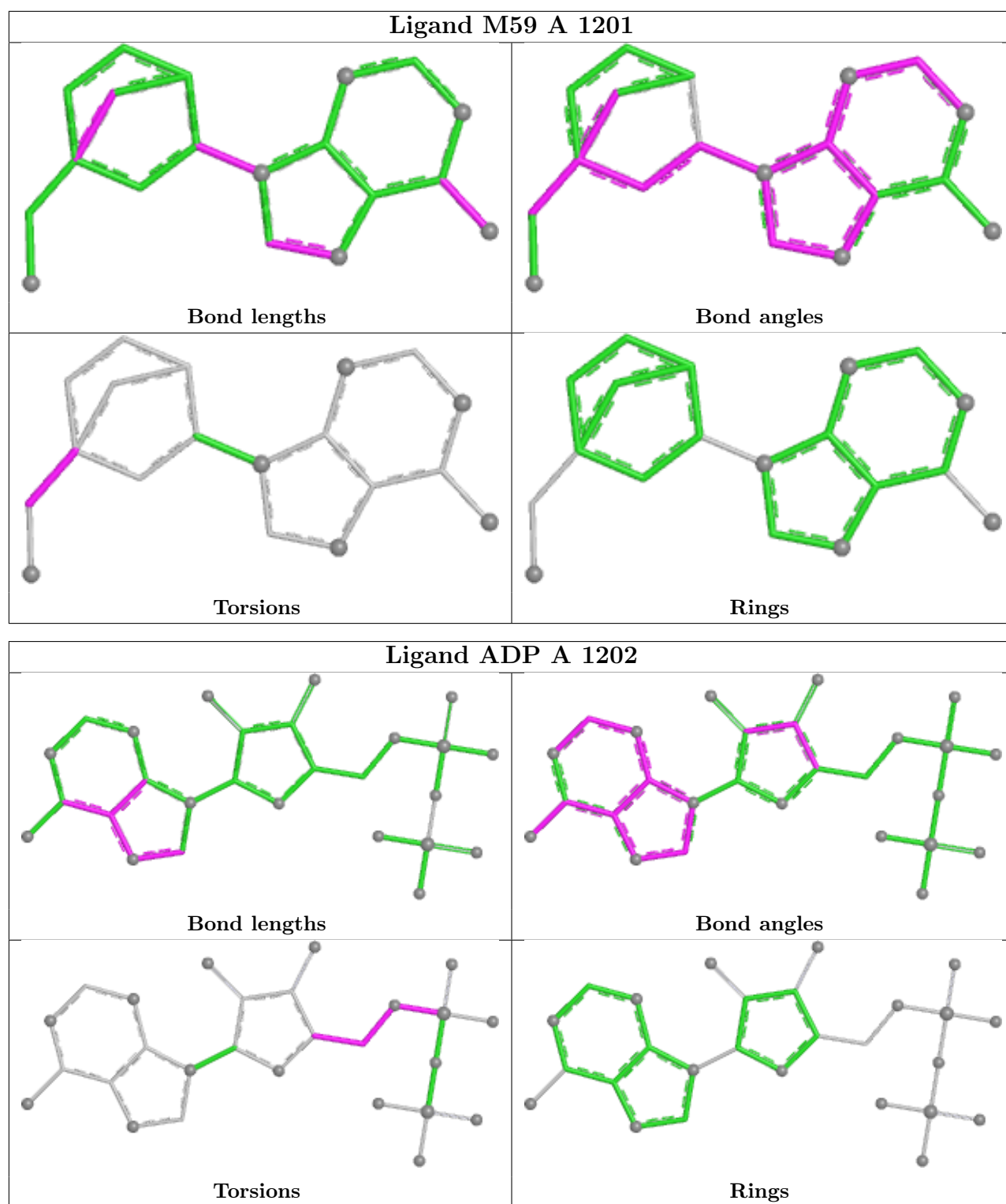
Mol	Chain	Res	Type	Atoms
2	A	1201	M59	O13-C12-C5-C3
2	A	1201	M59	O13-C12-C5-C8
2	A	1201	M59	O13-C12-C5-C11
3	A	1202	ADP	C5'-O5'-PA-O2A
3	A	1202	ADP	C5'-O5'-PA-O3A
3	A	1202	ADP	C4'-C5'-O5'-PA
3	A	1202	ADP	C3'-C4'-C5'-O5'
3	A	1202	ADP	O4'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	M59	3	0
3	A	1202	ADP	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 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	486/556 (87%)	0.54	31 (6%) 25 21	41, 58, 85, 104	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	332	TRP	5.4
1	A	1166	LEU	5.0
1	A	1165	ASN	4.3
1	A	181	ARG	3.9
1	A	183	CYS	3.8
1	A	1096	VAL	3.5
1	A	331	ASP	3.3
1	A	135	SER	3.3
1	A	454	ALA	3.3
1	A	166	TRP	3.2
1	A	134	SER	3.2
1	A	101	ARG	3.2
1	A	167	THR	3.1
1	A	169	TRP	3.0
1	A	182	ASP	2.9
1	A	186	LEU	2.7
1	A	132	GLN	2.6
1	A	131	TYR	2.6
1	A	333	VAL	2.5
1	A	1117	THR	2.4
1	A	159	TYR	2.3
1	A	232	ARG	2.2
1	A	1141	TYR	2.2
1	A	136	GLY	2.1
1	A	129	ARG	2.1
1	A	168	LYS	2.1
1	A	161	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	180	GLY	2.0
1	A	319	ASP	2.0
1	A	1159	THR	2.0
1	A	356	PRO	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](#)

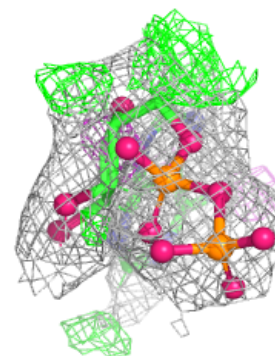
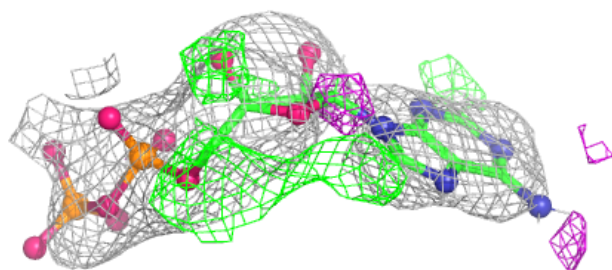
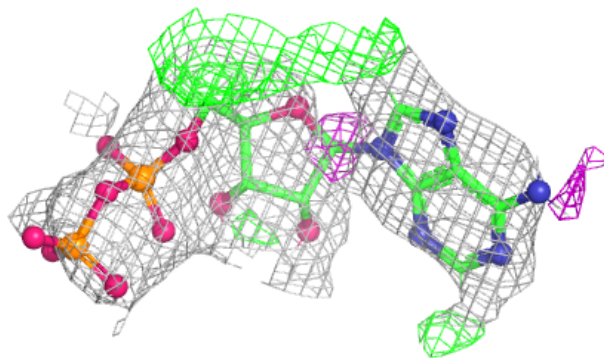
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ADP	A	1202	27/27	0.64	0.19	56,74,101,109	39
2	M59	A	1201	19/19	0.80	0.17	46,66,82,82	0

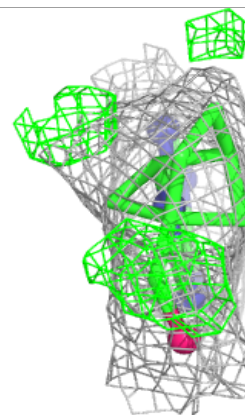
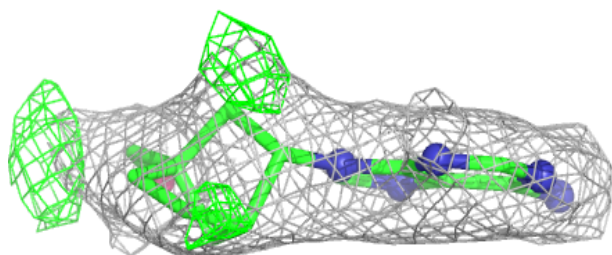
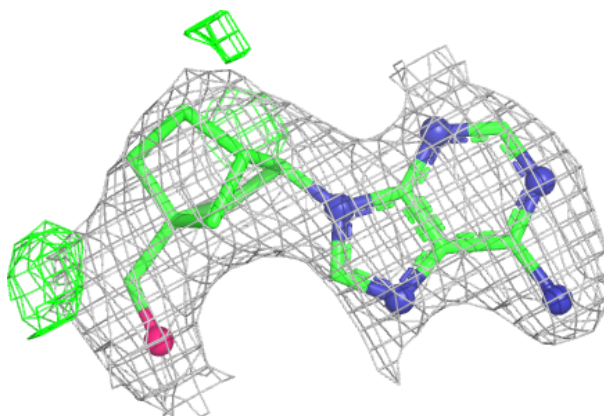
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 1202:

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

**Electron density around M59 A 1201:**

$2mF_o-DF_c$ (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.