



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:21 AM UTC

PDB ID : 2BKK / pdb_00002bkk
Title : Crystal structure of Aminoglycoside Phosphotransferase APH(3')-IIIa in complex with the inhibitor AR_3a
Authors : Kohl, A.; Amstutz, P.; Parizek, P.; Binz, H.K.; Briand, C.; Capitani, G.; Forrer, P.; Pluckthun, A.; Grutter, M.G.
Deposited on : 2005-02-16
Resolution : 2.15 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

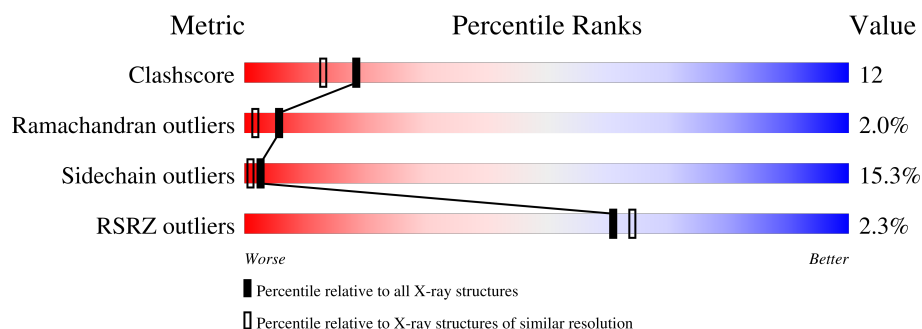
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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))
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	264	
1	C	264	
2	B	169	
2	D	169	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7000 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 AMINOGLYCOSIDE 3'-PHOSPHOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			2038	1304	326	398	10			
1	C	262	Total	C	N	O	S	0	0	0
			2165	1382	343	430	10			

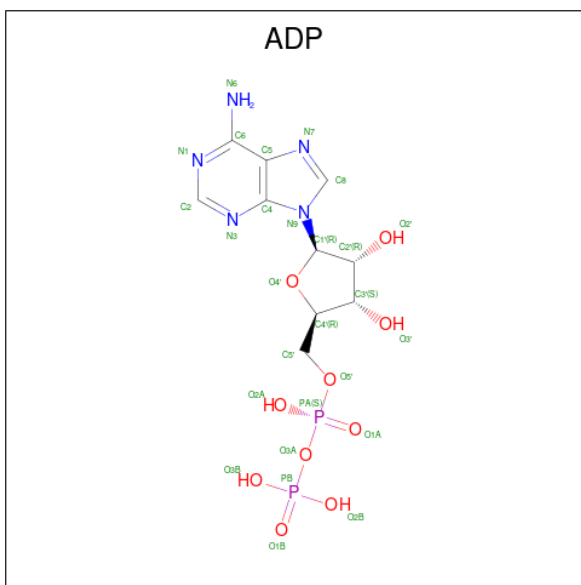
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	SER	CYS	engineered mutation	UNP P00554
C	19	SER	CYS	engineered mutation	UNP P00554

- Molecule 2 is a protein called DESIGNED ANKYRIN REPEAT INHIBITOR AR_3A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	156	Total	C	N	O	S	0	0	1
			1194	756	201	235	2			
2	D	157	Total	C	N	O	S	0	0	1
			1198	758	202	236	2			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
3	C	1	Total 27	C 10	N 5	O 10	P 2	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Mg 2 2	0	0
4	C	2	Total Mg 2 2	0	0

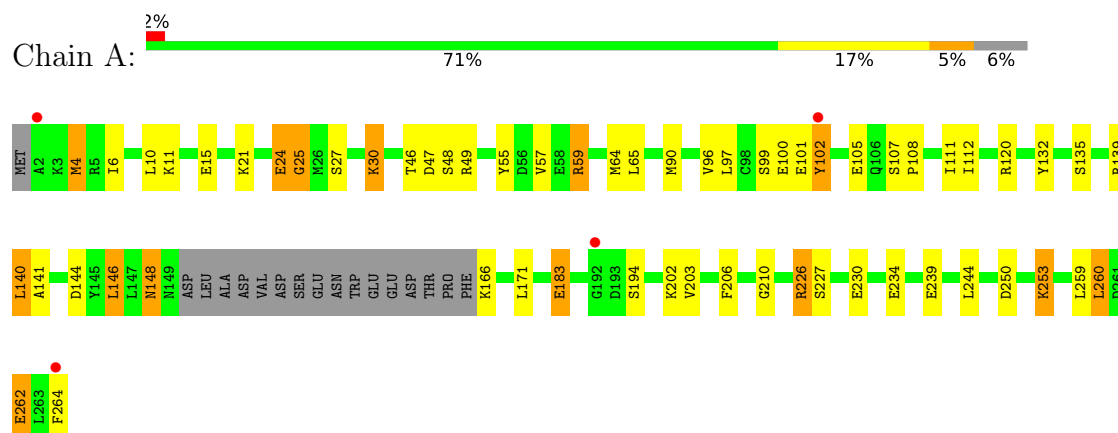
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	115	Total O 115 115	0	0
5	B	53	Total O 53 53	0	0
5	C	130	Total O 130 130	0	0
5	D	49	Total O 49 49	0	0

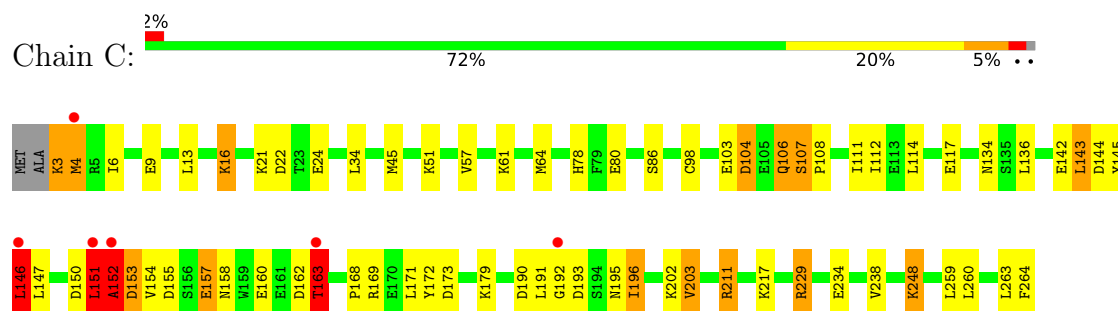
3 Residue-property plots

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.

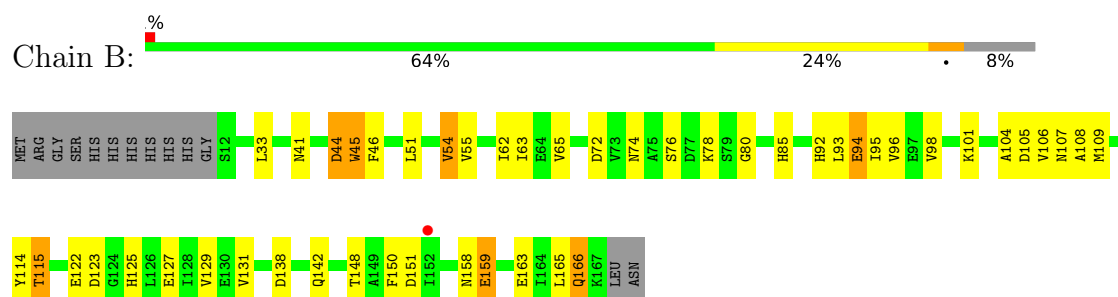
• Molecule 1: AMINOGLYCOSIDE 3'-PHOSPHOTRANSFERASE



• Molecule 1: AMINOGLYCOSIDE 3'-PHOSPHOTRANSFERASE

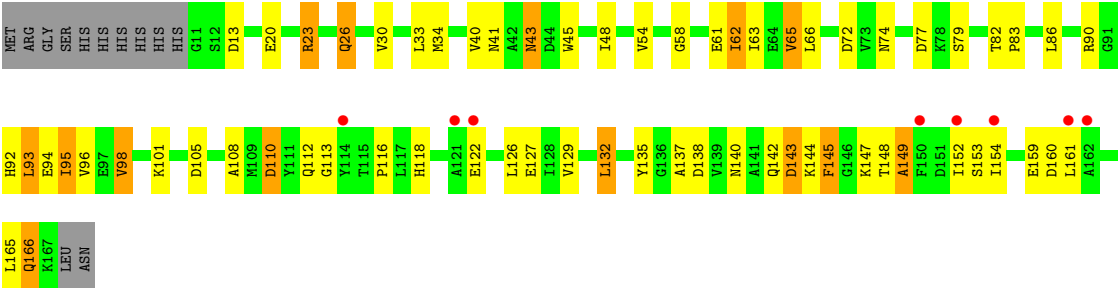


• Molecule 2: DESIGNED ANKYRIN REPEAT INHIBITOR AR_3A



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.67Å 98.08Å 81.30Å 90.00° 110.01° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15 20.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	96.7 (20.00-2.15) 96.9 (20.00-2.15)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.15Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.199 , 0.260 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7000	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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: MG, 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	0.96	2/2081 (0.1%)	0.98	0/2800
1	C	0.95	2/2213 (0.1%)	0.99	2/2984 (0.1%)
2	B	0.82	1/1215 (0.1%)	0.98	2/1652 (0.1%)
2	D	0.90	1/1219 (0.1%)	1.04	2/1657 (0.1%)
All	All	0.92	6/6728 (0.1%)	1.00	6/9093 (0.1%)

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

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	64	MET	SD-CE	-13.27	1.46	1.79
1	C	64	MET	SD-CE	-8.78	1.57	1.79
2	B	166	GLN	C-N	-6.86	1.23	1.33
2	D	166	GLN	C-N	-6.53	1.24	1.33
1	A	90	MET	SD-CE	5.68	1.93	1.79
1	C	146	LEU	N-CA	5.55	1.53	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	44	ASP	N-CA-C	-6.37	99.91	110.17
2	B	115	THR	N-CA-C	-6.14	101.00	110.58
2	D	98	VAL	N-CA-CB	5.60	116.73	110.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	45	TRP	CB-CA-C	-5.59	104.10	111.82
1	C	152	ALA	N-CA-C	-5.50	99.09	110.80
1	C	196	ILE	N-CA-C	5.35	115.59	108.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	152	ALA	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	2038	0	1992	41	0
1	C	2165	0	2091	41	0
2	B	1194	0	1162	33	0
2	D	1198	0	1165	46	0
3	A	27	0	12	1	0
3	C	27	0	12	0	0
4	A	2	0	0	0	0
4	C	2	0	0	0	0
5	A	115	0	0	12	1
5	B	53	0	0	10	0
5	C	130	0	0	8	1
5	D	49	0	0	13	0
All	All	7000	0	6434	157	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 12.

All (157) 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:C:146:LEU:HA	5:C:2081:HOH:O	1.34	1.23

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:CYS:SG	5:C:2066:HOH:O	1.96	1.19
2:D:82:THR:OG1	5:D:2030:HOH:O	1.74	1.04
2:D:116:PRO:HD3	5:D:2039:HOH:O	1.59	1.02
2:D:41:ASN:HD21	2:D:72:ASP:H	1.12	0.96
2:B:44:ASP:O	5:B:2025:HOH:O	1.84	0.93
2:D:43:ASN:HA	5:D:2019:HOH:O	1.69	0.90
2:D:132:LEU:O	5:D:2041:HOH:O	1.89	0.89
2:D:101:LYS:NZ	2:D:135:TYR:OH	2.07	0.87
2:B:74:ASN:HD21	2:B:105:ASP:H	1.23	0.86
2:D:108:ALA:O	5:D:2039:HOH:O	1.94	0.85
2:B:45:TRP:N	5:B:2026:HOH:O	2.05	0.81
2:D:74:ASN:HD21	2:D:105:ASP:H	1.27	0.79
1:A:47:ASP:OD1	5:A:2031:HOH:O	2.05	0.74
2:B:127:GLU:O	5:B:2045:HOH:O	2.04	0.73
2:D:48:ILE:O	5:D:2019:HOH:O	2.07	0.73
1:C:146:LEU:N	5:C:2082:HOH:O	2.22	0.72
1:C:190:ASP:O	1:C:195:ASN:ND2	2.21	0.72
1:A:141:ALA:HB2	5:B:2003:HOH:O	1.91	0.71
2:D:110:ASP:OD1	2:D:110:ASP:C	2.33	0.71
1:A:65:LEU:O	5:A:2037:HOH:O	2.10	0.70
2:B:106:VAL:O	5:B:2043:HOH:O	2.09	0.70
1:C:104:ASP:HB2	1:C:106:GLN:HE21	1.56	0.70
2:B:62:ILE:HD12	2:B:65:VAL:HG11	1.75	0.68
2:B:123:ASP:HB3	2:B:125:HIS:ND1	2.07	0.68
1:C:143:LEU:O	1:C:146:LEU:HB3	1.93	0.68
1:A:141:ALA:CB	5:B:2003:HOH:O	2.40	0.67
2:D:41:ASN:ND2	2:D:72:ASP:H	1.90	0.67
1:A:194:SER:OG	5:A:2085:HOH:O	2.12	0.66
2:B:108:ALA:O	2:B:115:THR:HA	1.95	0.66
1:A:97:LEU:HD21	5:A:2113:HOH:O	1.95	0.65
1:A:59:ARG:HD2	1:A:210:GLY:O	1.96	0.65
2:B:107:ASN:HD21	2:B:138:ASP:H	1.45	0.65
2:D:74:ASN:ND2	2:D:105:ASP:H	1.94	0.64
2:D:137:ALA:N	5:D:2041:HOH:O	2.32	0.63
2:D:43:ASN:HB2	5:D:2017:HOH:O	1.99	0.63
1:C:143:LEU:CD1	1:C:264:PHE:HB3	2.29	0.63
1:C:248:LYS:H	1:C:248:LYS:NZ	1.97	0.62
1:A:46:THR:HG23	1:A:57:VAL:HG13	1.81	0.62
1:C:191:LEU:HD12	1:C:196:ILE:HD11	1.80	0.62
1:A:46:THR:HG21	1:A:55:TYR:O	2.00	0.62
2:B:74:ASN:ND2	2:B:105:ASP:H	1.94	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:ALA:O	1:C:153:ASP:O	2.18	0.61
2:B:62:ILE:O	2:B:65:VAL:HG12	2.01	0.60
2:D:127:GLU:H	2:D:127:GLU:CD	2.09	0.60
1:A:97:LEU:CD2	5:A:2113:HOH:O	2.49	0.60
2:B:62:ILE:HA	2:B:65:VAL:HG12	1.83	0.60
2:D:41:ASN:HD21	2:D:72:ASP:N	1.94	0.60
2:D:126:LEU:O	2:D:129:VAL:HG12	2.01	0.60
1:C:144:ASP:O	1:C:147:LEU:HG	2.04	0.58
1:C:3:LYS:O	1:C:4:MET:O	2.21	0.58
1:A:140:LEU:HB3	2:B:45:TRP:CG	2.38	0.58
1:C:155:ASP:OD2	1:C:157:GLU:O	2.22	0.57
2:D:40:VAL:HG23	2:D:66:LEU:HD22	1.86	0.57
2:B:127:GLU:C	5:B:2045:HOH:O	2.46	0.57
1:C:248:LYS:H	1:C:248:LYS:HZ3	1.53	0.57
1:A:120:ARG:NH1	5:A:2064:HOH:O	2.38	0.56
1:A:97:LEU:HD11	1:A:194:SER:C	2.30	0.56
2:D:62:ILE:HA	2:D:65:VAL:HG13	1.86	0.55
1:C:106:GLN:OE1	1:C:107:SER:OG	2.24	0.55
2:D:149:ALA:N	5:D:2046:HOH:O	2.39	0.55
2:B:123:ASP:HB3	2:B:125:HIS:CE1	2.42	0.54
1:C:104:ASP:HB2	1:C:106:GLN:NE2	2.22	0.54
1:C:108:PRO:O	1:C:112:ILE:HG12	2.07	0.54
1:C:172:TYR:OH	2:D:48:ILE:HD12	2.09	0.53
2:D:74:ASN:HD21	2:D:105:ASP:N	2.02	0.53
2:D:77:ASP:C	2:D:77:ASP:OD1	2.52	0.52
2:B:62:ILE:O	2:B:65:VAL:CG1	2.57	0.52
1:A:250:ASP:CG	1:A:253:LYS:HD3	2.33	0.52
1:C:117:GLU:CD	1:C:203:VAL:HG12	2.35	0.52
1:A:6:ILE:CD1	1:A:10:LEU:HD23	2.41	0.51
1:A:105:GLU:O	1:A:105:GLU:HG3	2.11	0.51
2:B:131:VAL:HG23	5:B:2045:HOH:O	2.10	0.51
1:A:135:SER:O	1:A:139:ARG:HG3	2.10	0.51
1:C:146:LEU:CA	5:C:2082:HOH:O	2.59	0.51
1:A:226:ARG:NE	5:A:2094:HOH:O	2.31	0.50
1:C:106:GLN:NE2	1:C:106:GLN:H	2.09	0.50
2:D:93:LEU:O	2:D:96:VAL:HG12	2.12	0.50
1:C:151:LEU:HB2	1:C:153:ASP:HA	1.94	0.50
1:C:146:LEU:HG	1:C:146:LEU:O	2.11	0.49
1:A:4:MET:HE1	1:A:11:LYS:HD3	1.93	0.49
1:A:6:ILE:HD12	1:A:10:LEU:HD23	1.94	0.49
2:B:54:VAL:CG1	2:B:63:ILE:HG13	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:134:ASN:ND2	1:C:211:ARG:O	2.44	0.48
2:B:109:MET:HA	2:B:114:TYR:O	2.13	0.48
1:A:108:PRO:O	1:A:112:ILE:HG12	2.13	0.48
1:C:168:PRO:HG3	2:D:79:SER:HB3	1.96	0.48
2:D:20:GLU:OE2	2:D:23:ARG:NH1	2.47	0.48
2:D:92:HIS:HB2	5:D:2032:HOH:O	2.13	0.48
1:A:97:LEU:HD21	1:A:194:SER:HA	1.94	0.48
2:B:74:ASN:HD21	2:B:105:ASP:N	2.01	0.48
1:C:162:ASP:O	1:C:163:THR:O	2.31	0.48
2:B:62:ILE:CD1	2:B:65:VAL:HG11	2.43	0.48
2:D:160:ASP:HA	5:D:2048:HOH:O	2.13	0.48
2:D:138:ASP:OD1	2:D:138:ASP:C	2.57	0.47
2:B:44:ASP:CA	5:B:2026:HOH:O	2.62	0.47
2:B:51:LEU:O	2:B:55:VAL:HG23	2.15	0.47
2:D:34:MET:HE1	2:D:40:VAL:HB	1.96	0.47
1:A:226:ARG:NH1	5:A:2094:HOH:O	2.47	0.47
2:D:82:THR:HB	2:D:83:PRO:HD2	1.97	0.47
2:B:159:GLU:C	2:B:163:GLU:HG2	2.40	0.47
1:C:211:ARG:HB3	5:C:2105:HOH:O	2.16	0.46
1:A:59:ARG:CD	1:A:210:GLY:O	2.64	0.46
1:C:16:LYS:H	1:C:16:LYS:CD	2.28	0.46
2:D:93:LEU:HA	2:D:96:VAL:HG12	1.97	0.46
1:A:141:ALA:O	1:A:144:ASP:HB3	2.14	0.46
1:A:30:LYS:N	1:A:30:LYS:HD2	2.30	0.46
1:A:96:VAL:HG11	1:A:101:GLU:OE2	2.16	0.46
1:C:192:GLY:O	1:C:196:ILE:HD12	2.16	0.46
2:D:26:GLN:O	2:D:30:VAL:HG23	2.16	0.46
2:B:76:SER:OG	2:B:80:GLY:HA2	2.16	0.45
1:C:143:LEU:HD13	1:C:264:PHE:HB3	1.97	0.45
2:B:41:ASN:HD21	2:B:72:ASP:H	1.65	0.45
2:D:118:HIS:NE2	2:D:148:THR:O	2.49	0.45
2:B:45:TRP:HB2	2:B:46:PHE:CD1	2.52	0.45
2:D:30:VAL:O	2:D:34:MET:HG2	2.17	0.45
1:A:144:ASP:HB2	2:B:45:TRP:HA	1.99	0.45
2:B:85:HIS:HE1	2:B:108:ALA:HB3	1.82	0.45
1:C:264:PHE:O	5:C:2126:HOH:O	2.21	0.45
1:C:151:LEU:C	1:C:153:ASP:N	2.75	0.44
2:D:54:VAL:HG12	2:D:63:ILE:CG1	2.47	0.44
1:C:61:LYS:HD2	1:C:61:LYS:C	2.43	0.44
1:A:24:GLU:O	1:A:25:GLY:O	2.35	0.44
1:A:148:ASN:C	1:A:148:ASN:HD22	2.26	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:VAL:HG11	1:A:206:PHE:CE1	2.52	0.44
1:C:150:ASP:O	1:C:151:LEU:C	2.61	0.44
1:A:183:GLU:HG3	5:A:2080:HOH:O	2.18	0.43
1:A:250:ASP:OD2	1:A:253:LYS:HD3	2.18	0.43
2:D:143:ASP:OD1	5:D:2044:HOH:O	2.21	0.43
1:C:229:ARG:HD2	1:C:238:VAL:HG21	2.01	0.43
2:D:113:GLY:HA2	5:D:2040:HOH:O	2.17	0.43
1:A:27:SER:OG	3:A:1265:ADP:O2B	2.27	0.43
2:D:101:LYS:CE	2:D:135:TYR:OH	2.67	0.43
2:B:46:PHE:HA	2:B:78:LYS:HD2	2.01	0.42
1:C:45:MET:HA	1:C:86:SER:O	2.19	0.42
1:C:146:LEU:HA	5:C:2082:HOH:O	2.17	0.42
2:D:145:PHE:N	2:D:145:PHE:CD2	2.87	0.42
1:A:203:VAL:HG11	1:A:206:PHE:CZ	2.53	0.42
2:D:58:GLY:O	2:D:95:ILE:HD11	2.19	0.42
2:D:54:VAL:HG12	2:D:63:ILE:HG13	2.01	0.42
1:A:99:SER:HB3	5:A:2084:HOH:O	2.20	0.42
1:C:78:HIS:HE1	1:C:80:GLU:OE2	2.02	0.42
1:A:59:ARG:HG2	1:A:132:TYR:CE1	2.55	0.41
1:A:262:GLU:HA	1:A:264:PHE:CE2	2.55	0.41
1:A:15:GLU:HA	5:A:2008:HOH:O	2.20	0.41
2:B:74:ASN:HD21	2:B:104:ALA:HA	1.86	0.41
2:D:40:VAL:HG22	2:D:40:VAL:O	2.21	0.41
2:D:145:PHE:N	2:D:145:PHE:HD2	2.19	0.41
1:A:55:TYR:CE1	1:A:146:LEU:HD11	2.56	0.41
2:B:94:GLU:O	2:B:98:VAL:HG23	2.20	0.41
1:C:104:ASP:CB	1:C:106:GLN:HE21	2.28	0.40
2:D:86:LEU:HD23	2:D:86:LEU:O	2.21	0.40
1:C:21:LYS:HB3	5:C:2015:HOH:O	2.20	0.40
1:A:260:LEU:O	5:A:2108:HOH:O	2.22	0.40
2:B:92:HIS:HB2	5:B:2039:HOH:O	2.22	0.40
1:C:169:ARG:O	1:C:173:ASP:OD1	2.38	0.40
2:D:48:ILE:N	2:D:48:ILE:HD13	2.36	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)
5:A:2109:HOH:O	5:C:2056:HOH:O[2_545]	1.89	0.31

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	243/264 (92%)	229 (94%)	12 (5%)	2 (1%)	16	10
1	C	260/264 (98%)	241 (93%)	10 (4%)	9 (4%)	3	0
2	B	154/169 (91%)	141 (92%)	11 (7%)	2 (1%)	9	5
2	D	155/169 (92%)	130 (84%)	22 (14%)	3 (2%)	6	2
All	All	812/866 (94%)	741 (91%)	55 (7%)	16 (2%)	6	2

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	GLY
1	C	4	MET
1	C	153	ASP
1	C	163	THR
1	C	193	ASP
1	A	102	TYR
2	B	148	THR
1	C	146	LEU
1	C	152	ALA
2	D	143	ASP
2	D	149	ALA
2	B	45	TRP
1	C	145	TYR
1	C	151	LEU
1	C	22	ASP
2	D	145	PHE

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	226/242 (93%)	198 (88%)	28 (12%)	4	2
1	C	241/242 (100%)	204 (85%)	37 (15%)	2	1
2	B	123/135 (91%)	107 (87%)	16 (13%)	4	1
2	D	123/135 (91%)	95 (77%)	28 (23%)	1	0
All	All	713/754 (95%)	604 (85%)	109 (15%)	3	1

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	MET
1	A	21	LYS
1	A	24	GLU
1	A	30	LYS
1	A	48	SER
1	A	49	ARG
1	A	59	ARG
1	A	100	GLU
1	A	102	TYR
1	A	107	SER
1	A	111	ILE
1	A	140	LEU
1	A	146	LEU
1	A	148	ASN
1	A	166	LYS
1	A	171	LEU
1	A	183	GLU
1	A	202	LYS
1	A	226	ARG
1	A	227	SER
1	A	230	GLU
1	A	234	GLU
1	A	239	GLU
1	A	244	LEU
1	A	253	LYS
1	A	259	LEU
1	A	260	LEU
1	A	262	GLU
2	B	33	LEU
2	B	54	VAL

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Mol	Chain	Res	Type
2	B	93	LEU
2	B	94	GLU
2	B	95	ILE
2	B	96	VAL
2	B	101	LYS
2	B	122	GLU
2	B	129	VAL
2	B	142	GLN
2	B	150	PHE
2	B	151	ASP
2	B	158	ASN
2	B	159	GLU
2	B	165	LEU
2	B	166	GLN
1	C	3	LYS
1	C	6	ILE
1	C	9	GLU
1	C	13	LEU
1	C	16	LYS
1	C	24	GLU
1	C	34	LEU
1	C	51	LYS
1	C	57	VAL
1	C	103	GLU
1	C	104	ASP
1	C	106	GLN
1	C	107	SER
1	C	111	ILE
1	C	114	LEU
1	C	136	LEU
1	C	142	GLU
1	C	143	LEU
1	C	146	LEU
1	C	151	LEU
1	C	154	VAL
1	C	157	GLU
1	C	158	ASN
1	C	160	GLU
1	C	163	THR
1	C	171	LEU
1	C	179	LYS
1	C	202	LYS

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Mol	Chain	Res	Type
1	C	203	VAL
1	C	211	ARG
1	C	217	LYS
1	C	229	ARG
1	C	234	GLU
1	C	248	LYS
1	C	259	LEU
1	C	260	LEU
1	C	263	LEU
2	D	13	ASP
2	D	23	ARG
2	D	26	GLN
2	D	33	LEU
2	D	43	ASN
2	D	61	GLU
2	D	62	ILE
2	D	65	VAL
2	D	90	ARG
2	D	93	LEU
2	D	94	GLU
2	D	95	ILE
2	D	98	VAL
2	D	110	ASP
2	D	112	GLN
2	D	122	GLU
2	D	132	LEU
2	D	140	ASN
2	D	142	GLN
2	D	144	LYS
2	D	147	LYS
2	D	152	ILE
2	D	153	SER
2	D	154	ILE
2	D	159	GLU
2	D	161	LEU
2	D	165	LEU
2	D	166	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	HIS

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Mol	Chain	Res	Type
1	A	148	ASN
1	A	236	GLN
2	B	26	GLN
2	B	36	ASN
2	B	41	ASN
2	B	43	ASN
2	B	59	HIS
2	B	74	ASN
2	B	107	ASN
2	B	118	HIS
2	B	158	ASN
1	C	78	HIS
1	C	148	ASN
2	D	36	ASN
2	D	41	ASN
2	D	56	ASN
2	D	74	ASN
2	D	140	ASN
2	D	156	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 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	$\# Z > 2$
3	ADP	A	1265	4	28,29,29	1.45	5 (17%)	43,45,45	1.69	10 (23%)
3	ADP	C	1265	4	28,29,29	1.12	2 (7%)	43,45,45	1.91	8 (18%)

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	ADP	A	1265	4	-	3/16/32/32	0/3/3/3
3	ADP	C	1265	4	-	0/16/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1265	ADP	C2-N1	3.68	1.40	1.33
3	C	1265	ADP	C2-N1	2.89	1.39	1.33
3	A	1265	ADP	C8-N7	2.73	1.36	1.31
3	A	1265	ADP	C2-N3	2.70	1.38	1.33
3	A	1265	ADP	PB-O1B	2.13	1.57	1.50
3	A	1265	ADP	PA-O3A	2.08	1.61	1.59
3	C	1265	ADP	PA-O3A	2.05	1.61	1.59

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1265	ADP	N3-C2-N1	-7.94	116.56	128.58
3	A	1265	ADP	N3-C2-N1	-4.82	121.28	128.58
3	A	1265	ADP	C5-C4-N3	-4.64	120.33	126.72
3	C	1265	ADP	C2-N1-C6	3.78	124.94	118.73
3	C	1265	ADP	C2-N3-C4	3.61	120.66	111.83
3	C	1265	ADP	C5-C4-N3	-3.32	122.15	126.72
3	A	1265	ADP	O2A-PA-O3A	3.24	116.02	107.27
3	A	1265	ADP	C2-N3-C4	2.97	119.08	111.83
3	C	1265	ADP	N9-C8-N7	-2.93	109.77	113.94
3	A	1265	ADP	N9-C8-N7	-2.55	110.33	113.94
3	A	1265	ADP	C5-N7-C8	2.54	107.44	103.45
3	A	1265	ADP	N3-C4-N9	2.52	131.45	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1265	ADP	C6-C5-C4	2.46	120.54	117.18
3	C	1265	ADP	C5-N7-C8	2.35	107.14	103.45
3	C	1265	ADP	O4'-C1'-N9	2.35	112.59	108.09
3	A	1265	ADP	C4-C5-N7	-2.29	107.97	110.58
3	A	1265	ADP	C5-C4-N9	2.19	108.20	105.81
3	C	1265	ADP	O2B-PB-O3A	2.13	111.79	104.64

There are no chirality outliers.

All (3) torsion outliers are listed below:

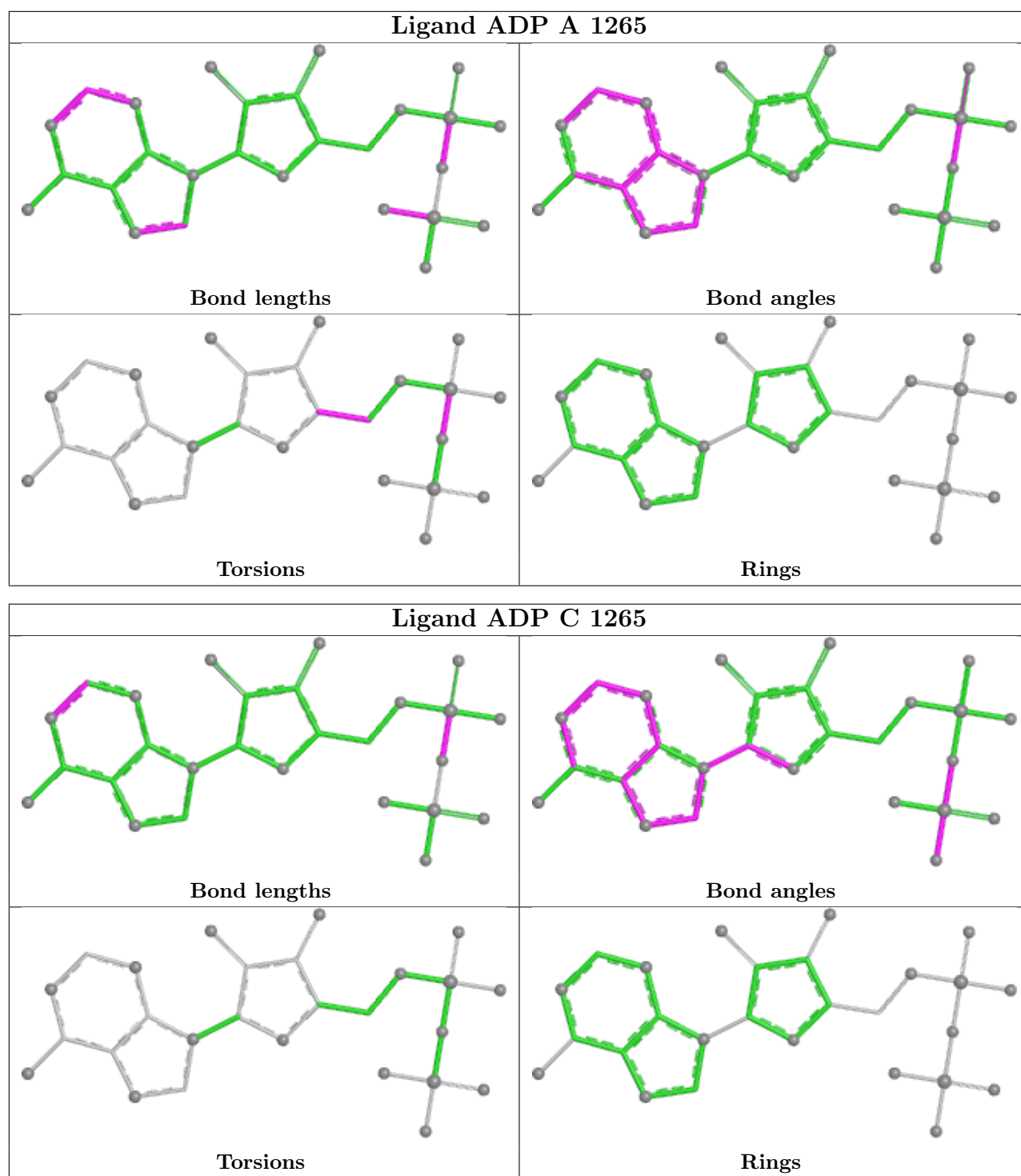
Mol	Chain	Res	Type	Atoms
3	A	1265	ADP	PB-O3A-PA-O2A
3	A	1265	ADP	O4'-C4'-C5'-O5'
3	A	1265	ADP	PB-O3A-PA-O1A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1265	ADP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

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	247/264 (93%)	0.02	4 (1%) 70 74	29, 50, 87, 115	0
1	C	262/264 (99%)	0.03	6 (2%) 61 64	29, 47, 83, 112	1 (0%)
2	B	156/169 (92%)	0.22	1 (0%) 85 87	34, 59, 117, 123	0
2	D	157/169 (92%)	0.50	8 (5%) 33 37	31, 59, 117, 128	0
All	All	822/866 (94%)	0.15	19 (2%) 61 64	29, 52, 105, 128	1 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	152	ALA	5.0
1	A	264	PHE	3.4
1	C	151	LEU	3.2
1	A	102	TYR	3.1
2	D	162	ALA	3.0
2	D	152	ILE	2.8
2	D	154	ILE	2.8
2	D	114	TYR	2.4
2	D	161	LEU	2.3
1	C	163	THR	2.2
2	B	152	ILE	2.2
1	C	146	LEU	2.2
2	D	122	GLU	2.1
1	A	2	ALA	2.1
2	D	121	ALA	2.1
1	A	192	GLY	2.0
1	C	192	GLY	2.0
2	D	150	PHE	2.0
1	C	4	MET	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](#)

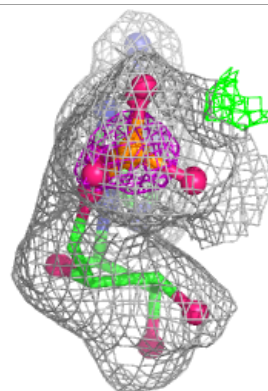
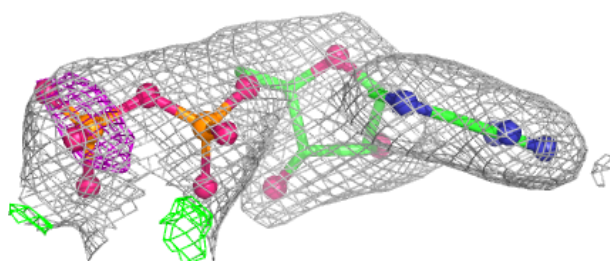
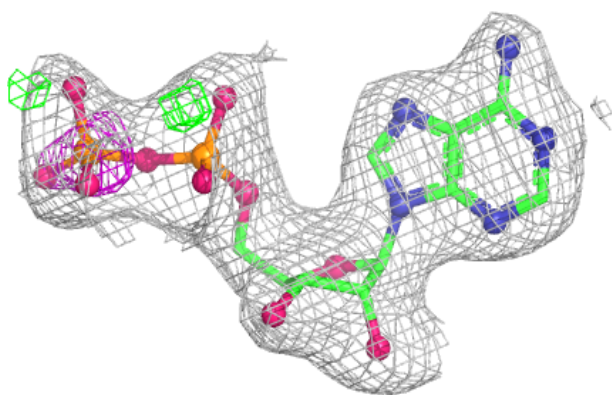
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	MG	C	1267	1/1	0.72	0.14	91,91,91,91	0
4	MG	C	1266	1/1	0.84	0.07	64,64,64,64	0
4	MG	A	1267	1/1	0.86	0.10	60,60,60,60	0
3	ADP	C	1265	27/27	0.93	0.07	26,42,71,74	0
3	ADP	A	1265	27/27	0.94	0.06	30,38,48,53	0
4	MG	A	1266	1/1	0.98	0.04	50,50,50,50	0

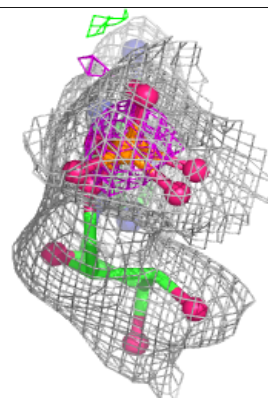
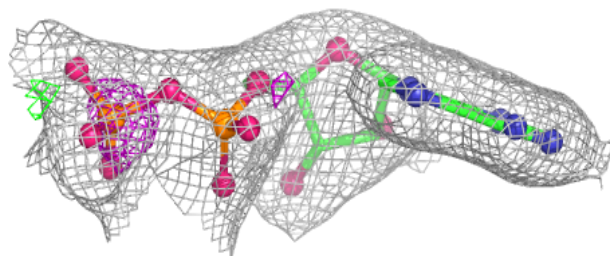
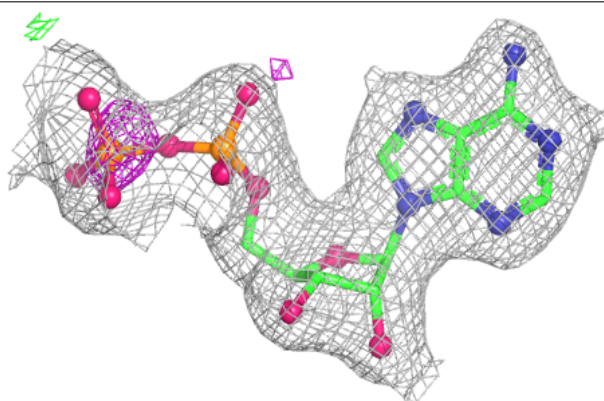
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP C 1265:

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

**Electron density around ADP A 1265:**

$2mF_o-DF_c$ (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.