



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:40 AM UTC

PDB ID : 2I8C / pdb_00002i8c
Title : Allosteric inhibition of Staphylococcus aureus D-alanine:D-alanine ligase revealed by crystallographic studies
Authors : Liu, S.; Chang, J.S.; Herberg, J.T.; Horng, M.; Tomich, P.K.; Lin, A.H.; Marotti, K.R.
Deposited on : 2006-09-01
Resolution : 2.46 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

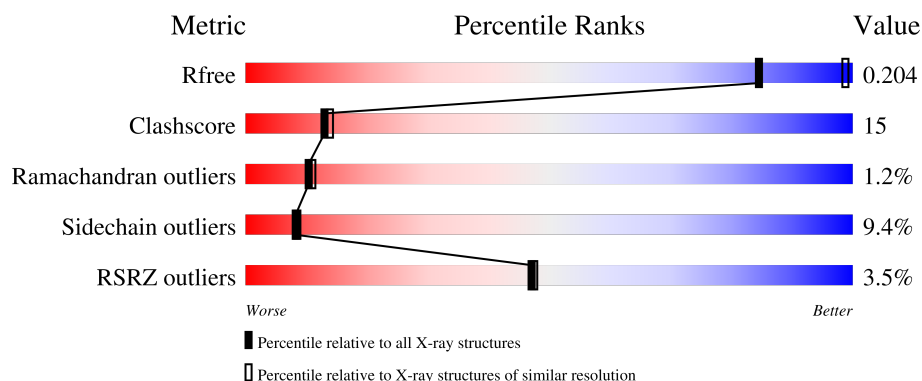
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	4% 69% 22% • 5%
1	B	364	2% 60% 31% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5561 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.

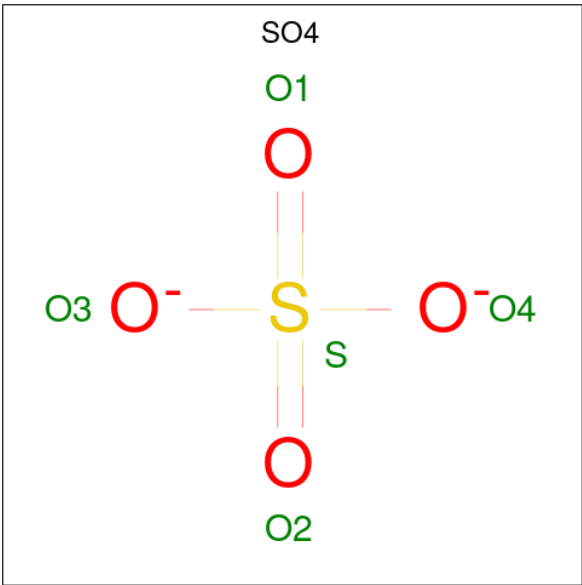
- Molecule 1 is a protein called D-alanine-D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2738	1741	453	535	9			
1	B	347	Total	C	N	O	S	0	0	0
			2758	1753	459	537	9			

There are 16 discrepancies between the modelled and reference sequences:

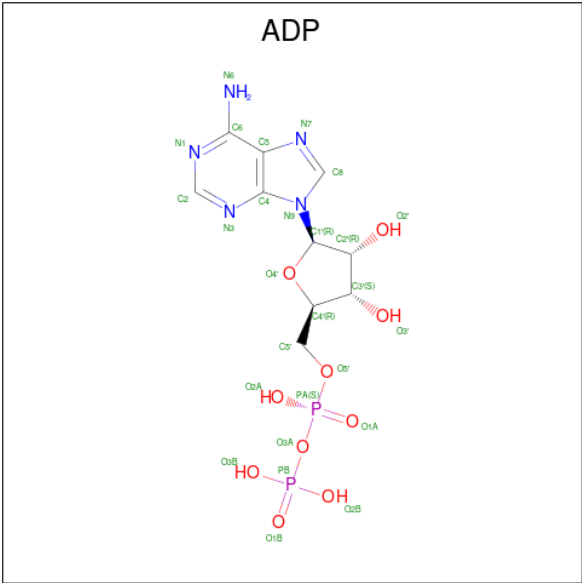
Chain	Residue	Modelled	Actual	Comment	Reference
A	357	ARG	-	cloning artifact	UNP Q5HEB7
A	358	SER	-	cloning artifact	UNP Q5HEB7
A	359	HIS	-	expression tag	UNP Q5HEB7
A	360	HIS	-	expression tag	UNP Q5HEB7
A	361	HIS	-	expression tag	UNP Q5HEB7
A	362	HIS	-	expression tag	UNP Q5HEB7
A	363	HIS	-	expression tag	UNP Q5HEB7
A	364	HIS	-	expression tag	UNP Q5HEB7
B	357	ARG	-	cloning artifact	UNP Q5HEB7
B	358	SER	-	cloning artifact	UNP Q5HEB7
B	359	HIS	-	expression tag	UNP Q5HEB7
B	360	HIS	-	expression tag	UNP Q5HEB7
B	361	HIS	-	expression tag	UNP Q5HEB7
B	362	HIS	-	expression tag	UNP Q5HEB7
B	363	HIS	-	expression tag	UNP Q5HEB7
B	364	HIS	-	expression tag	UNP Q5HEB7

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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



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

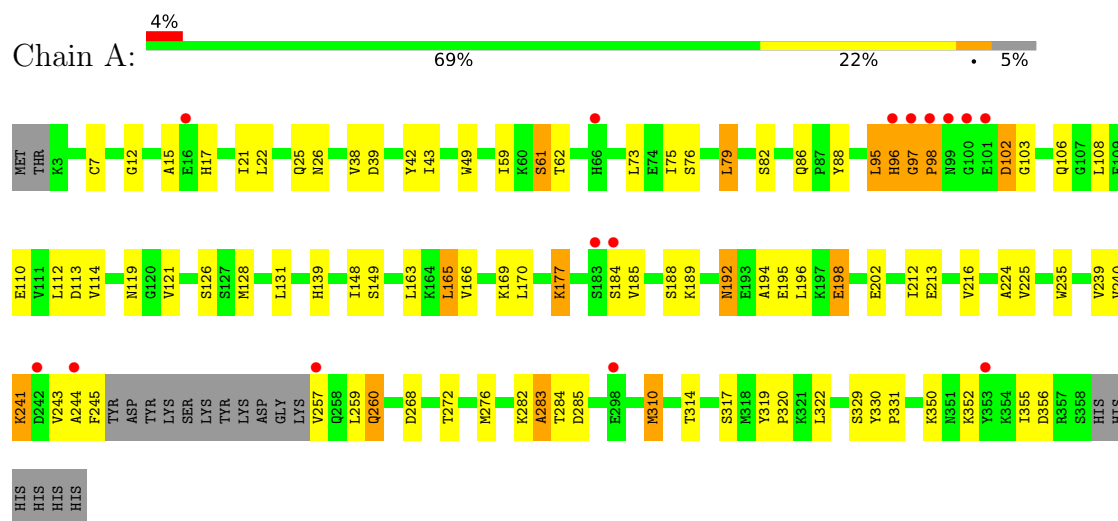
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		

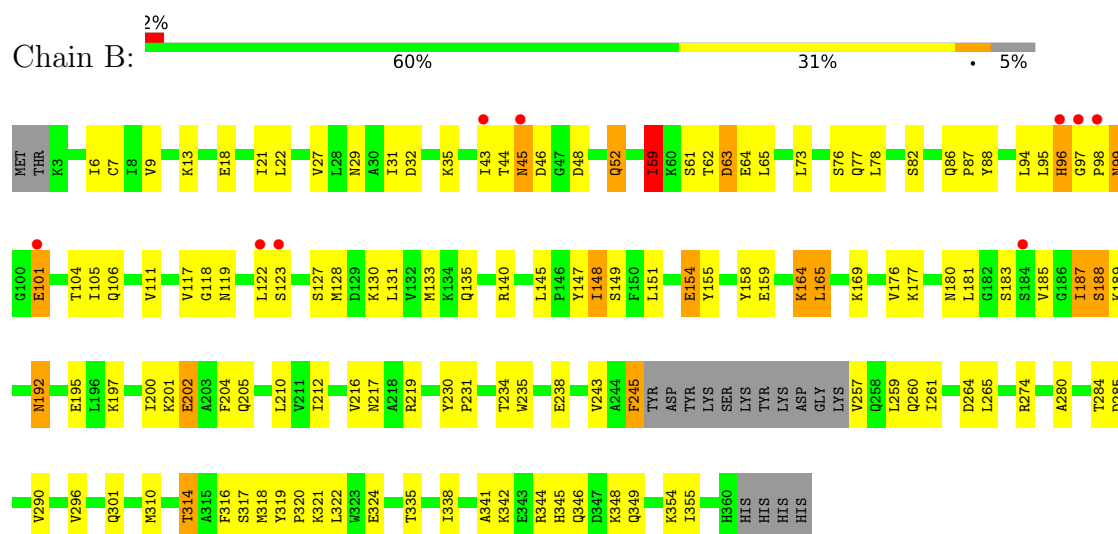
3 Residue-property plots [i](#)

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 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: D-alanine-D-alanine ligase



• Molecule 1: D-alanine-D-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.80Å 66.97Å 78.99Å 90.00° 96.05° 90.00°	Depositor
Resolution (Å)	47.51 – 2.46 47.51 – 2.46	Depositor EDS
% Data completeness (in resolution range)	97.2 (47.51-2.46) 97.4 (47.51-2.46)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.3.0008	Depositor
R, R_{free}	0.210 , 0.269 0.217 , 0.204	Depositor DCC
R_{free} test set	1289 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.114	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.5	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	0.93	EDS
Total number of atoms	5561	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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: ADP, MG, SO4

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	1.10	4/2790 (0.1%)	1.12	6/3775 (0.2%)
1	B	1.08	2/2812 (0.1%)	1.08	6/3805 (0.2%)
All	All	1.09	6/5602 (0.1%)	1.10	12/7580 (0.2%)

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 (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	VAL	CA-CB	6.38	1.62	1.54
1	A	21	ILE	CA-CB	5.44	1.61	1.54
1	A	283	ALA	CA-CB	-5.35	1.45	1.53
1	B	118	GLY	C-O	-5.17	1.18	1.23
1	A	213	GLU	CA-C	-5.12	1.46	1.52
1	B	280	ALA	C-O	-5.08	1.18	1.24

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	177	LYS	CA-C-N	-6.28	113.51	119.85
1	B	177	LYS	C-N-CA	-6.28	113.51	119.85
1	B	338	ILE	CB-CA-C	-6.05	104.22	111.97
1	A	97	GLY	CA-C-N	5.75	127.03	119.84
1	A	97	GLY	C-N-CA	5.75	127.03	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	GLY	N-CA-C	-5.62	107.22	113.79
1	B	338	ILE	N-CA-C	5.59	115.78	110.42
1	A	114	VAL	CB-CA-C	-5.43	104.77	110.53
1	A	61	SER	N-CA-C	5.27	115.31	107.88
1	B	119	ASN	N-CA-C	5.18	117.36	110.53
1	B	78	LEU	N-CA-C	5.17	118.69	112.38
1	A	62	THR	N-CA-C	5.02	117.64	111.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	184	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	2738	0	2700	67	0
1	B	2758	0	2714	104	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	27	0	12	1	0
3	B	27	0	12	0	0
4	B	1	0	0	0	0
All	All	5561	0	5438	165	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 15.

All (165) 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:102:ASP:HB2	1:A:128:MET:HE3	1.37	1.02
1:B:95:LEU:C	1:B:310:MET:HE1	1.93	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ILE:HD12	1:A:165:LEU:HD13	1.53	0.90
1:B:127:SER:HA	1:B:133:MET:HE3	1.62	0.81
1:B:101:GLU:HB3	1:B:128:MET:HE2	1.62	0.81
1:B:131:LEU:HD11	1:B:135:GLN:NE2	1.97	0.80
1:A:43:ILE:HD12	1:A:95:LEU:HD11	1.66	0.76
1:B:145:LEU:HD21	1:B:216:VAL:HG11	1.70	0.73
1:B:181:LEU:HD22	1:B:181:LEU:N	2.03	0.73
1:B:314:THR:CG2	1:B:316:PHE:H	2.02	0.72
1:B:131:LEU:HD11	1:B:135:GLN:HE21	1.55	0.71
1:B:96:HIS:N	1:B:310:MET:HE1	2.05	0.71
1:B:13:LYS:HA	1:B:45:ASN:HA	1.73	0.71
1:A:76:SER:H	1:A:79:LEU:HD23	1.56	0.70
1:A:43:ILE:HD12	1:A:95:LEU:CD1	2.22	0.69
1:A:76:SER:O	1:A:79:LEU:HD23	1.93	0.69
1:A:25:GLN:HE21	1:A:59:ILE:HG22	1.58	0.67
1:A:244:ALA:O	1:A:245:PHE:O	2.15	0.65
1:B:164:LYS:CE	1:B:165:LEU:HD23	2.28	0.64
1:B:314:THR:HG22	1:B:316:PHE:N	2.13	0.64
1:B:127:SER:CA	1:B:133:MET:HE3	2.28	0.63
1:A:73:LEU:HD21	1:A:79:LEU:HD13	1.80	0.62
1:B:59:ILE:HD11	1:B:65:LEU:HD11	1.82	0.61
1:B:127:SER:HA	1:B:133:MET:CE	2.30	0.61
1:A:76:SER:H	1:A:79:LEU:CD2	2.13	0.60
1:A:148:ILE:CD1	1:A:165:LEU:HD13	2.28	0.60
1:B:314:THR:HB	1:B:317:SER:OG	2.01	0.60
1:B:131:LEU:CD1	1:B:135:GLN:NE2	2.65	0.59
1:B:314:THR:HG23	1:B:316:PHE:H	1.67	0.59
1:A:102:ASP:HB2	1:A:128:MET:CE	2.22	0.59
1:A:96:HIS:O	1:A:310:MET:HE2	2.02	0.59
1:B:145:LEU:HD21	1:B:216:VAL:CG1	2.32	0.59
1:B:101:GLU:O	1:B:104:THR:HG23	2.03	0.59
1:A:243:VAL:HG12	1:A:245:PHE:H	1.67	0.58
1:A:163:LEU:HD22	1:A:196:LEU:HD23	1.85	0.58
1:B:245:PHE:O	1:B:245:PHE:HD1	1.86	0.58
1:B:181:LEU:N	1:B:181:LEU:CD2	2.67	0.57
1:B:219:ARG:NH2	1:B:265:LEU:HD23	2.20	0.57
1:A:352:LYS:O	1:A:356:ASP:OD1	2.22	0.56
1:A:49:TRP:CD2	1:A:75:ILE:HD11	2.40	0.56
1:B:86:GLN:HB2	1:B:87:PRO:HD2	1.88	0.56
1:A:26:ASN:ND2	1:A:314:THR:HB	2.21	0.56
1:A:49:TRP:CE2	1:A:75:ILE:HD11	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:ASN:ND2	1:B:195:GLU:H	2.05	0.55
1:B:122:LEU:O	1:B:122:LEU:HD12	2.08	0.53
1:B:13:LYS:H	1:B:98:PRO:HG3	1.72	0.53
1:B:97:GLY:O	1:B:101:GLU:HA	2.08	0.53
1:A:7:CYS:HA	1:A:39:ASP:O	2.09	0.52
1:B:43:ILE:HD12	1:B:95:LEU:CD1	2.39	0.52
1:B:147:TYR:O	1:B:148:ILE:HD12	2.09	0.52
1:A:17:HIS:NE2	1:A:42:TYR:OH	2.24	0.52
1:B:101:GLU:HB3	1:B:128:MET:CE	2.36	0.52
1:A:240:VAL:HG22	1:A:260:GLN:HB2	1.91	0.52
1:A:106:GLN:HE22	1:A:310:MET:HE1	1.75	0.51
1:B:147:TYR:O	1:B:148:ILE:CD1	2.58	0.51
1:A:177:LYS:NZ	3:A:400:ADP:N7	2.55	0.51
1:A:82:SER:HB3	1:A:86:GLN:O	2.10	0.51
1:A:239:VAL:O	1:A:241:LYS:HG3	2.11	0.51
1:B:18:GLU:HA	1:B:21:ILE:HD12	1.92	0.51
1:B:151:LEU:HB2	1:B:154:GLU:HG3	1.93	0.50
1:B:219:ARG:NH2	1:B:265:LEU:CD2	2.74	0.50
1:A:76:SER:N	1:A:79:LEU:HD23	2.24	0.50
1:A:97:GLY:O	1:A:98:PRO:O	2.29	0.50
1:A:73:LEU:HD21	1:A:79:LEU:CD1	2.42	0.50
1:B:165:LEU:HD13	1:B:169:LYS:HD2	1.94	0.49
1:B:345:HIS:CE1	1:B:349:GLN:NE2	2.80	0.49
1:A:192:ASN:ND2	1:A:195:GLU:H	2.10	0.49
1:B:59:ILE:CD1	1:B:65:LEU:HD11	2.42	0.49
1:B:95:LEU:CA	1:B:310:MET:HE1	2.41	0.49
1:B:95:LEU:O	1:B:95:LEU:HG	2.13	0.49
1:B:44:THR:HG22	1:B:48:ASP:HB3	1.95	0.49
1:A:7:CYS:HB2	1:A:88:TYR:CE2	2.47	0.48
1:A:139:HIS:CG	1:B:135:GLN:HG2	2.47	0.48
1:A:192:ASN:ND2	1:A:192:ASN:C	2.71	0.48
1:A:49:TRP:CE3	1:A:75:ILE:HD11	2.48	0.48
1:B:234:THR:O	1:B:274:ARG:HG2	2.13	0.48
1:B:43:ILE:HD12	1:B:95:LEU:HD11	1.94	0.48
1:A:192:ASN:C	1:A:192:ASN:HD22	2.21	0.48
1:B:260:GLN:NE2	1:B:264:ASP:H	2.11	0.48
1:B:314:THR:HG22	1:B:317:SER:N	2.28	0.48
1:B:319:TYR:HB3	1:B:320:PRO:HD3	1.95	0.48
1:B:13:LYS:HG2	1:B:45:ASN:HA	1.96	0.48
1:A:26:ASN:HD22	1:A:314:THR:HB	1.80	0.47
1:B:59:ILE:HD11	1:B:65:LEU:CD1	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:LEU:N	1:B:310:MET:HE1	2.29	0.47
1:B:245:PHE:CD1	1:B:245:PHE:C	2.92	0.47
1:A:25:GLN:HE22	1:A:61:SER:HA	1.79	0.47
1:B:245:PHE:HD1	1:B:245:PHE:C	2.23	0.47
1:A:196:LEU:HD11	1:A:212:ILE:HD12	1.97	0.46
1:A:192:ASN:HD21	1:A:194:ALA:HB3	1.81	0.46
1:A:317:SER:HB2	1:A:320:PRO:HD2	1.97	0.46
1:B:176:VAL:HG22	1:B:212:ILE:HD12	1.97	0.46
1:A:75:ILE:HB	1:B:76:SER:HB2	1.97	0.46
1:A:198:GLU:O	1:A:202:GLU:HG3	2.15	0.46
1:B:59:ILE:HD12	1:B:65:LEU:HD21	1.97	0.46
1:B:123:SER:OG	1:B:284:THR:O	2.28	0.46
1:B:158:TYR:O	1:B:159:GLU:C	2.58	0.46
1:B:131:LEU:HD13	1:B:149:SER:HB3	1.98	0.46
1:A:272:THR:HG22	1:A:276:MET:HE3	1.97	0.45
1:B:127:SER:C	1:B:133:MET:HE3	2.41	0.45
1:B:32:ASP:HB3	1:B:35:LYS:HD3	1.99	0.45
1:A:319:TYR:HB3	1:A:320:PRO:HD3	1.98	0.45
1:B:44:THR:HG23	1:B:46:ASP:N	2.32	0.45
1:B:155:TYR:CZ	1:B:200:ILE:HG21	2.52	0.45
1:B:106:GLN:H	1:B:106:GLN:CD	2.25	0.45
1:B:202:GLU:OE1	1:B:202:GLU:CA	2.65	0.45
1:A:355:ILE:CD1	1:B:181:LEU:HD11	2.47	0.44
1:B:205:GLN:HE21	1:B:205:GLN:HA	1.82	0.44
1:B:342:LYS:HE2	1:B:346:GLN:HE22	1.82	0.44
1:B:344:ARG:O	1:B:348:LYS:HG2	2.17	0.44
1:A:235:TRP:O	1:A:322:LEU:HD22	2.18	0.44
1:B:164:LYS:NZ	1:B:165:LEU:HD23	2.32	0.44
1:B:318:MET:HG3	1:B:322:LEU:HD12	1.99	0.44
1:A:108:LEU:C	1:A:108:LEU:HD23	2.42	0.44
1:B:219:ARG:HD2	1:B:238:GLU:OE1	2.18	0.44
1:B:318:MET:HG3	1:B:322:LEU:CD1	2.48	0.44
1:B:63:ASP:N	1:B:63:ASP:OD1	2.49	0.44
1:B:235:TRP:O	1:B:322:LEU:HD22	2.18	0.43
1:B:127:SER:CA	1:B:133:MET:CE	2.94	0.43
1:B:29:ASN:OD1	1:B:59:ILE:HG23	2.18	0.43
1:B:131:LEU:HD13	1:B:149:SER:CB	2.49	0.43
1:A:330:TYR:HB3	1:A:331:PRO:HD3	2.00	0.43
1:B:202:GLU:OE1	1:B:202:GLU:HA	2.17	0.43
1:A:329:SER:OG	1:A:331:PRO:HD2	2.18	0.43
1:B:140:ARG:O	1:B:140:ARG:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:LEU:HD12	1:B:111:VAL:HG11	2.01	0.42
1:B:6:ILE:HG22	1:B:7:CYS:N	2.34	0.42
1:A:131:LEU:HD22	1:A:149:SER:HB2	2.01	0.42
1:B:181:LEU:HD13	1:B:181:LEU:HA	1.79	0.42
1:B:192:ASN:HD22	1:B:195:GLU:H	1.66	0.42
1:B:44:THR:CG2	1:B:48:ASP:HB3	2.50	0.42
1:B:155:TYR:CE2	1:B:200:ILE:HG21	2.54	0.42
1:A:282:LYS:O	1:A:283:ALA:C	2.61	0.42
1:B:96:HIS:N	1:B:310:MET:CE	2.80	0.42
1:B:230:TYR:N	1:B:231:PRO:CD	2.83	0.42
1:A:75:ILE:HA	1:A:79:LEU:HD21	2.02	0.42
1:A:15:ALA:C	1:A:17:HIS:H	2.27	0.42
1:B:187:ILE:HG22	1:B:188:SER:N	2.33	0.42
1:A:166:VAL:O	1:A:170:LEU:HB2	2.20	0.42
1:A:185:VAL:HG21	1:B:355:ILE:HG12	2.02	0.42
1:B:180:ASN:C	1:B:181:LEU:HD22	2.45	0.42
1:B:261:ILE:HD12	1:B:321:LYS:HB3	2.02	0.42
1:A:224:ALA:HB1	1:A:319:TYR:CE1	2.55	0.41
1:A:49:TRP:CZ2	1:A:75:ILE:HD11	2.55	0.41
1:A:49:TRP:CZ3	1:A:75:ILE:HD11	2.55	0.41
1:B:27:VAL:HG21	1:B:94:LEU:HD11	2.01	0.41
1:B:117:VAL:HA	1:B:341:ALA:HB2	2.01	0.41
1:A:112:LEU:O	1:A:113:ASP:HB3	2.20	0.41
1:B:82:SER:HA	1:B:88:TYR:CZ	2.55	0.41
1:A:12:GLY:O	1:A:17:HIS:ND1	2.52	0.41
1:A:165:LEU:HD23	1:A:165:LEU:HA	1.91	0.41
1:A:272:THR:CG2	1:A:276:MET:HE3	2.51	0.41
1:B:9:VAL:O	1:B:95:LEU:HD22	2.21	0.41
1:B:145:LEU:CD2	1:B:216:VAL:CG1	2.96	0.41
1:A:106:GLN:HG2	1:A:119:ASN:O	2.21	0.41
1:B:52:GLN:HE21	1:B:52:GLN:HB2	1.70	0.41
1:B:73:LEU:HB3	1:B:77:GLN:HB2	2.03	0.41
1:B:59:ILE:O	1:B:59:ILE:HG12	2.20	0.40
1:B:61:SER:OG	1:B:62:THR:N	2.53	0.40
1:A:76:SER:N	1:A:79:LEU:CD2	2.82	0.40
1:B:204:PHE:CE2	1:B:210:LEU:HG	2.56	0.40
1:B:32:ASP:CB	1:B:35:LYS:HD3	2.51	0.40
1:B:128:MET:HE3	1:B:128:MET:HB3	1.95	0.40
1:A:110:GLU:O	1:B:99:ASN:HB3	2.21	0.40
1:A:112:LEU:O	1:A:113:ASP:CB	2.69	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	341/364 (94%)	322 (94%)	17 (5%)	2 (1%)	21	27
1	B	343/364 (94%)	320 (93%)	17 (5%)	6 (2%)	7	6
All	All	684/728 (94%)	642 (94%)	34 (5%)	8 (1%)	10	11

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	HIS
1	A	98	PRO
1	B	59	ILE
1	B	99	ASN
1	B	96	HIS
1	B	101	GLU
1	B	183	SER
1	B	185	VAL

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	301/319 (94%)	277 (92%)	24 (8%)	11	13
1	B	303/319 (95%)	270 (89%)	33 (11%)	6	5
All	All	604/638 (95%)	547 (91%)	57 (9%)	8	8

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	38	VAL
1	A	79	LEU
1	A	95	LEU
1	A	102	ASP
1	A	126	SER
1	A	165	LEU
1	A	169	LYS
1	A	177	LYS
1	A	188	SER
1	A	189	LYS
1	A	192	ASN
1	A	198	GLU
1	A	216	VAL
1	A	225	VAL
1	A	241	LYS
1	A	257	VAL
1	A	259	LEU
1	A	260	GLN
1	A	268	ASP
1	A	284	THR
1	A	285	ASP
1	A	310	MET
1	A	350	LYS
1	B	22	LEU
1	B	31	ILE
1	B	45	ASN
1	B	52	GLN
1	B	59	ILE
1	B	63	ASP
1	B	64	GLU
1	B	105	ILE
1	B	130	LYS
1	B	148	ILE
1	B	154	GLU
1	B	164	LYS
1	B	165	LEU
1	B	187	ILE
1	B	188	SER
1	B	189	LYS
1	B	192	ASN
1	B	197	LYS
1	B	201	LYS

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Mol	Chain	Res	Type
1	B	202	GLU
1	B	217	ASN
1	B	243	VAL
1	B	245	PHE
1	B	257	VAL
1	B	259	LEU
1	B	285	ASP
1	B	290	VAL
1	B	296	VAL
1	B	301	GLN
1	B	314	THR
1	B	324	GLU
1	B	335	THR
1	B	354	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	86	GLN
1	A	135	GLN
1	A	161	ASN
1	A	192	ASN
1	A	345	HIS
1	B	26	ASN
1	B	45	ASN
1	B	52	GLN
1	B	77	GLN
1	B	99	ASN
1	B	135	GLN
1	B	161	ASN
1	B	192	ASN
1	B	205	GLN
1	B	260	GLN
1	B	301	GLN
1	B	345	HIS
1	B	346	GLN
1	B	349	GLN
1	B	359	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	400	-	28,29,29	1.41	5 (17%)	43,45,45	2.11	11 (25%)
2	SO4	B	401	-	4,4,4	0.23	0	6,6,6	0.69	0
3	ADP	B	400	4	28,29,29	1.68	5 (17%)	43,45,45	2.28	11 (25%)
2	SO4	A	401	-	4,4,4	0.27	0	6,6,6	0.20	0

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	400	-	-	6/16/32/32	0/3/3/3
3	ADP	B	400	4	-	5/16/32/32	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	400	ADP	C5-C4	5.38	1.48	1.39
3	A	400	ADP	C5-C4	4.50	1.47	1.39
3	B	400	ADP	PA-O3A	3.04	1.62	1.59
3	B	400	ADP	C5-N7	-2.94	1.33	1.39
3	B	400	ADP	C8-N7	2.71	1.36	1.31
3	A	400	ADP	PA-O3A	2.65	1.62	1.59
3	B	400	ADP	C5-C6	2.56	1.48	1.41
3	A	400	ADP	C5-C6	2.46	1.47	1.41
3	A	400	ADP	C5-N7	-2.34	1.34	1.39
3	A	400	ADP	C8-N7	2.22	1.35	1.31

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	400	ADP	C5-C4-N3	-7.78	116.00	126.72
3	B	400	ADP	N3-C4-N9	5.70	136.86	127.17
3	A	400	ADP	C5-C4-N3	-5.67	118.90	126.72
3	A	400	ADP	N3-C4-N9	4.69	135.15	127.17
3	B	400	ADP	C2-N3-C4	4.32	122.39	111.83
3	A	400	ADP	C2'-C1'-N9	-4.29	102.66	113.30
3	B	400	ADP	C4-C5-N7	-3.81	106.23	110.58
3	A	400	ADP	C4-C5-N7	-3.63	106.43	110.58
3	A	400	ADP	C2-N3-C4	3.54	120.48	111.83
3	B	400	ADP	C2'-C1'-N9	-3.34	105.01	113.30
3	A	400	ADP	O4'-C1'-N9	3.33	114.47	108.09
3	A	400	ADP	C4-N9-C8	3.31	109.21	105.74
3	B	400	ADP	C3'-C2'-C1'	3.25	107.61	101.46
3	A	400	ADP	N3-C2-N1	-3.15	123.81	128.58
3	B	400	ADP	N3-C2-N1	-3.05	123.97	128.58
3	A	400	ADP	C5-N7-C8	3.01	108.18	103.45
3	B	400	ADP	C5-N7-C8	2.92	108.03	103.45
3	A	400	ADP	O2A-PA-O3A	2.68	114.51	107.27
3	B	400	ADP	O4'-C4'-C5'	2.64	117.79	109.33
3	A	400	ADP	N9-C8-N7	-2.62	110.22	113.94
3	B	400	ADP	O5'-C5'-C4'	2.36	117.03	108.99
3	B	400	ADP	C1'-N9-C8	-2.04	122.58	127.09

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	400	ADP	PA-O3A-PB-O2B

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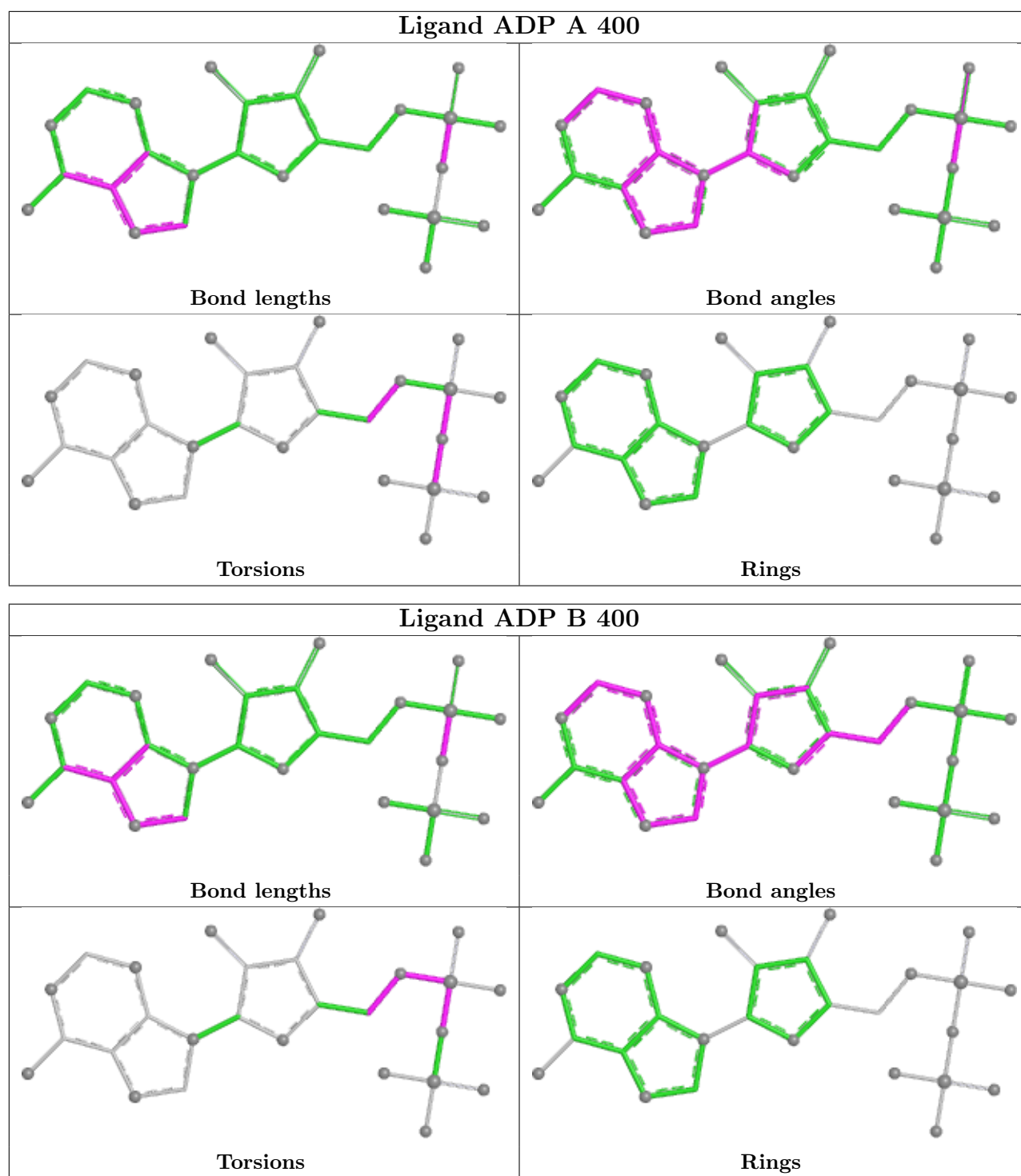
Mol	Chain	Res	Type	Atoms
3	B	400	ADP	PB-O3A-PA-O5'
3	B	400	ADP	C5'-O5'-PA-O2A
3	B	400	ADP	C4'-C5'-O5'-PA
3	B	400	ADP	C5'-O5'-PA-O1A
3	B	400	ADP	C5'-O5'-PA-O3A
3	A	400	ADP	PB-O3A-PA-O1A
3	A	400	ADP	PB-O3A-PA-O2A
3	A	400	ADP	PA-O3A-PB-O1B
3	A	400	ADP	PA-O3A-PB-O3B
3	A	400	ADP	C4'-C5'-O5'-PA

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	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	345/364 (94%)	0.40	15 (4%) 40 39	38, 52, 69, 81	0
1	B	347/364 (95%)	0.48	9 (2%) 57 58	37, 54, 68, 79	0
All	All	692/728 (95%)	0.44	24 (3%) 47 47	37, 53, 69, 81	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	GLY	5.7
1	A	183	SER	4.2
1	B	96	HIS	4.2
1	A	97	GLY	3.8
1	B	98	PRO	3.8
1	A	257	VAL	3.7
1	B	122	LEU	3.5
1	B	97	GLY	3.3
1	A	244	ALA	2.8
1	A	242	ASP	2.8
1	B	45	ASN	2.8
1	B	101	GLU	2.8
1	B	123	SER	2.7
1	A	16	GLU	2.7
1	A	99	ASN	2.6
1	A	96	HIS	2.4
1	A	101	GLU	2.4
1	A	98	PRO	2.3
1	B	43	ILE	2.2
1	A	66	HIS	2.2
1	A	353	TYR	2.2
1	A	298	GLU	2.1
1	A	184	SER	2.0
1	B	184	SER	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](#)

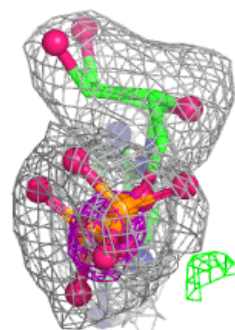
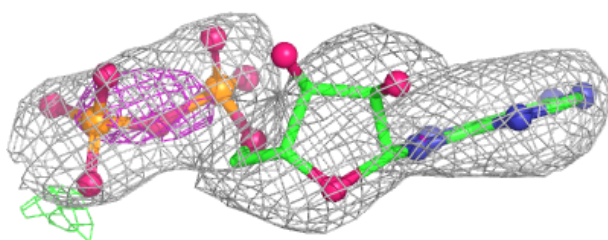
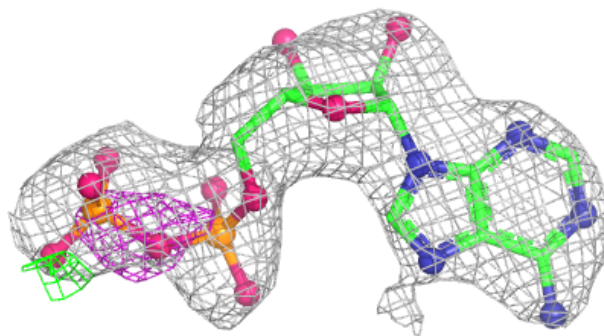
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B	402	1/1	0.80	0.10	80,80,80,80	0
3	ADP	A	400	27/27	0.84	0.13	88,89,93,94	0
2	SO4	A	401	5/5	0.85	0.10	92,93,94,94	0
2	SO4	B	401	5/5	0.85	0.15	83,84,85,85	0
3	ADP	B	400	27/27	0.89	0.10	53,60,73,75	0

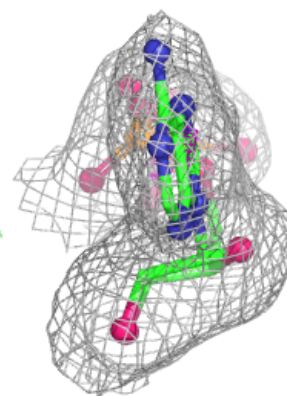
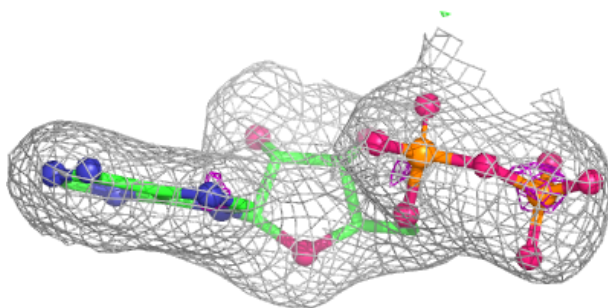
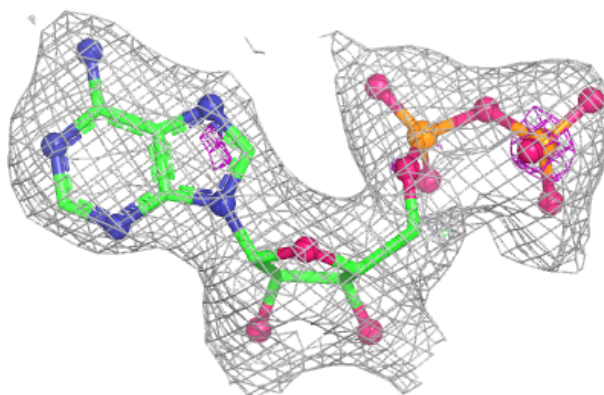
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 400:

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

**Electron density around ADP B 400:**

$2mF_o-DF_c$ (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.