



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:06 PM UTC

PDB ID : 4DHF / pdb_00004dhf
Title : Structure of Aurora A mutant bound to Biogenidec cpd 15
Authors : Silvian, L.; Marcotte, D.J.
Deposited on : 2012-01-27
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 ⓘ 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

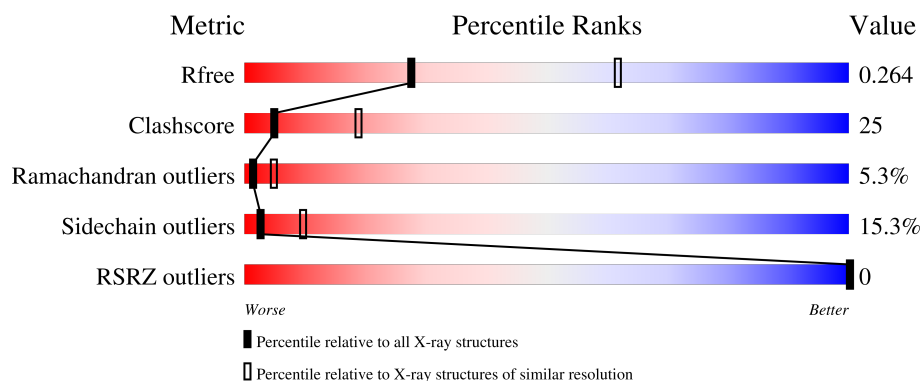
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 $\leq 5\%$. 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	271	
1	B	271	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4300 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.

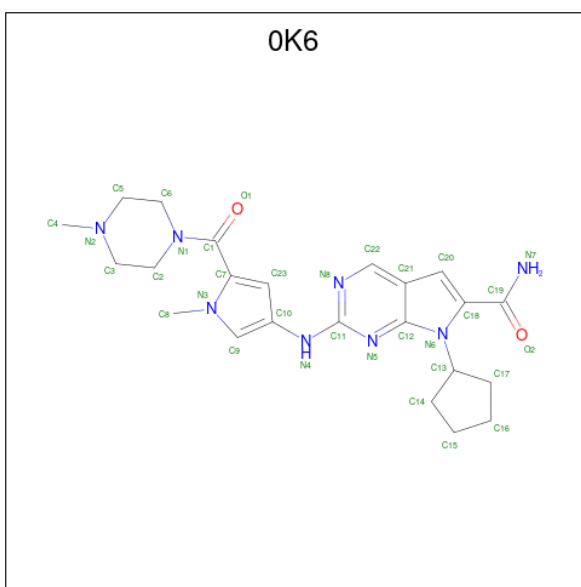
- Molecule 1 is a protein called Aurora kinase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			2111	1358	370	377	6			
1	B	258	Total	C	N	O	S	0	0	0
			2111	1358	370	377	6			

There are 14 discrepancies between the modelled and reference sequences:

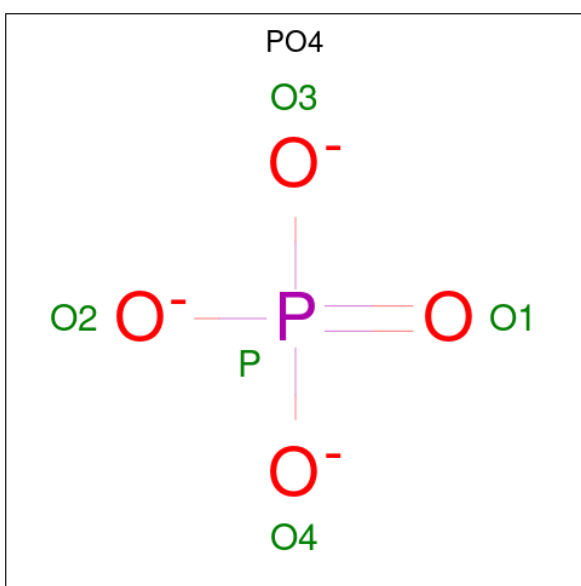
Chain	Residue	Modelled	Actual	Comment	Reference
A	121	GLY	-	expression tag	UNP O14965
A	122	PRO	-	expression tag	UNP O14965
A	123	LEU	-	expression tag	UNP O14965
A	124	GLY	-	expression tag	UNP O14965
A	125	SER	-	expression tag	UNP O14965
A	287	ASP	THR	engineered mutation	UNP O14965
A	288	ASP	THR	engineered mutation	UNP O14965
B	121	GLY	-	expression tag	UNP O14965
B	122	PRO	-	expression tag	UNP O14965
B	123	LEU	-	expression tag	UNP O14965
B	124	GLY	-	expression tag	UNP O14965
B	125	SER	-	expression tag	UNP O14965
B	287	ASP	THR	engineered mutation	UNP O14965
B	288	ASP	THR	engineered mutation	UNP O14965

- Molecule 2 is 7-cyclopentyl-2-({1-methyl-5-[(4-methylpiperazin-1-yl)carbonyl]-1H-pyrrol-3-yl}amino)-7H-pyrrolo[2,3-d]pyrimidine-6-carboxamide (CCD ID: 0K6) (formula: C₂₃H₃₀N₈O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	23	8	2		
2	B	1	Total	C	N	O	0	0
			33	23	8	2		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

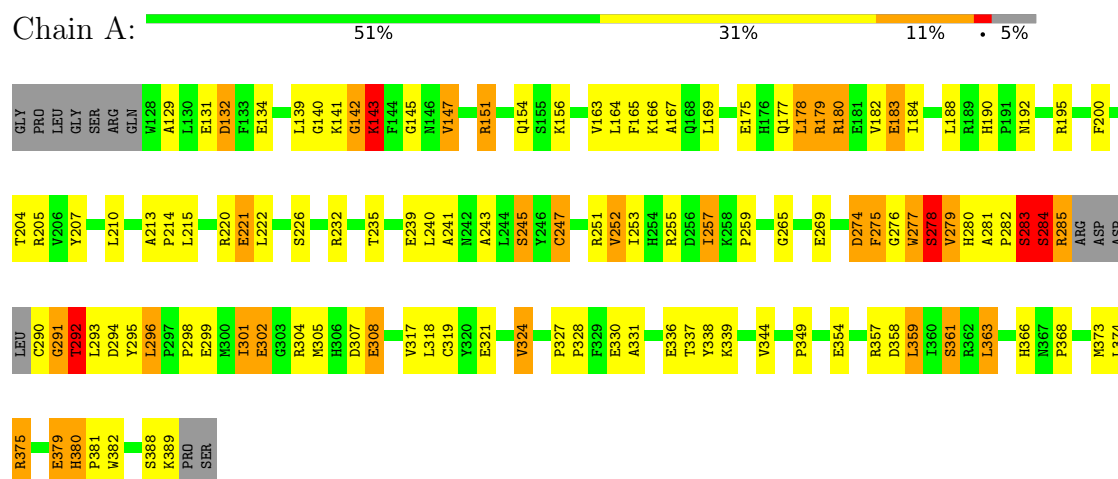
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

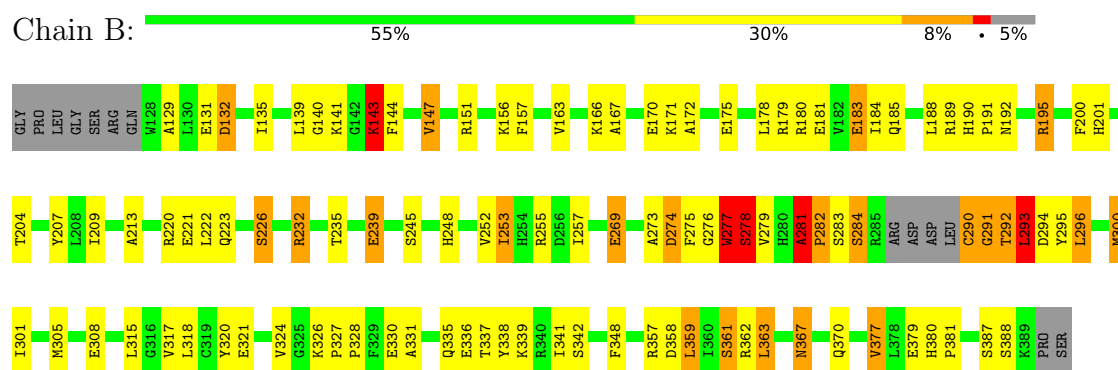
3 Residue-property plots

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: Aurora kinase A



• Molecule 1: Aurora kinase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	81.59Å 81.59Å 166.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.80 27.68 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.80) 99.7 (27.68-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	(Not available) , (Not available) (Not available) , 0.264	Depositor DCC
R_{free} test set	766 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	65.8	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.489 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4300	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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, PO4, 0K6

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	0.89	0/2164	1.08	5/2925 (0.2%)
1	B	0.90	0/2164	1.12	11/2925 (0.4%)
All	All	0.89	0/4328	1.10	16/5850 (0.3%)

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 (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	348	PHE	CA-C-N	-6.95	113.15	120.03
1	B	348	PHE	C-N-CA	-6.95	113.15	120.03
1	B	282	PRO	N-CA-C	6.39	121.94	113.57
1	B	213	ALA	CA-C-N	6.25	126.16	119.28
1	B	213	ALA	C-N-CA	6.25	126.16	119.28
1	B	278	SER	N-CA-C	6.00	127.81	111.00
1	B	294	ASP	N-CA-C	5.90	119.31	111.75
1	A	283	SER	CA-C-N	5.79	132.12	121.70
1	A	283	SER	C-N-CA	5.79	132.12	121.70
1	B	277	TRP	CA-C-N	5.75	132.05	121.70
1	B	277	TRP	C-N-CA	5.75	132.05	121.70
1	B	281	ALA	CA-C-N	5.34	124.85	119.19
1	B	281	ALA	C-N-CA	5.34	124.85	119.19
1	A	240	LEU	N-CA-C	5.13	116.56	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	324	VAL	CB-CA-C	-5.11	106.56	111.06
1	A	182	VAL	N-CA-C	5.01	115.13	110.42

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	278	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	2111	0	2107	113	0
1	B	2111	0	2107	106	0
2	A	33	0	30	9	0
2	B	33	0	30	6	0
3	A	5	0	0	1	0
3	B	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	4300	0	4274	216	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 25.

All (216) 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:183:GLU:CD	1:A:183:GLU:H	1.45	1.17
1:B:283:SER:HA	1:B:284:SER:HB3	1.28	1.15
1:A:375:ARG:H	1:A:375:ARG:HD2	1.15	1.12
1:B:183:GLU:N	1:B:183:GLU:CD	2.17	0.99
1:A:183:GLU:CD	1:A:183:GLU:N	2.14	0.98
1:B:183:GLU:CD	1:B:183:GLU:H	1.72	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:GLU:OE2	1:B:336:GLU:OE1	1.85	0.95
1:A:190:HIS:HD2	1:A:192:ASN:H	1.12	0.93
1:A:339:LYS:HZ2	1:B:339:LYS:HZ1	1.01	0.93
1:B:283:SER:HA	1:B:284:SER:CB	2.00	0.91
1:A:339:LYS:NZ	1:B:339:LYS:HZ1	1.67	0.91
1:A:283:SER:OG	1:A:284:SER:HB2	1.69	0.91
1:B:220:ARG:NH1	1:B:223:GLN:HB3	1.85	0.90
1:A:339:LYS:HZ2	1:B:339:LYS:NZ	1.68	0.90
1:A:375:ARG:HD2	1:A:375:ARG:N	1.91	0.86
1:B:140:GLY:HA3	2:B:401:OK6:H6	1.57	0.86
1:B:277:TRP:HA	1:B:278:SER:O	1.75	0.86
1:B:277:TRP:CD1	1:B:277:TRP:H	1.90	0.84
1:B:281:ALA:HB1	1:B:282:PRO:CD	2.08	0.83
1:A:283:SER:CB	1:A:284:SER:HB2	2.10	0.82
1:B:147:VAL:HG11	2:B:401:OK6:H2	1.43	0.81
1:B:184:ILE:O	1:B:188:LEU:HG	1.80	0.81
1:A:129:ALA:O	1:A:132:ASP:HB2	1.80	0.80
1:B:195:ARG:HH11	1:B:195:ARG:HG2	1.45	0.80
1:B:129:ALA:O	1:B:132:ASP:HB2	1.82	0.80
1:A:301:ILE:HG23	1:A:338:TYR:CE2	2.18	0.79
1:B:367:ASN:ND2	1:B:370:GLN:HG3	1.98	0.78
1:A:183:GLU:N	1:A:183:GLU:OE1	2.17	0.78
1:B:301:ILE:HG23	1:B:338:TYR:CE2	2.19	0.78
1:A:339:LYS:NZ	1:B:339:LYS:NZ	2.30	0.76
1:B:220:ARG:HH12	1:B:223:GLN:HB3	1.48	0.76
1:A:252:VAL:HB	1:A:279:VAL:HG11	1.67	0.76
1:B:283:SER:CA	1:B:284:SER:HB3	2.12	0.76
1:A:147:VAL:HG11	2:A:401:OK6:H2	1.53	0.73
1:B:232:ARG:HG3	1:B:232:ARG:HH11	1.54	0.73
1:A:291:GLY:O	1:A:293:LEU:N	2.21	0.73
1:B:281:ALA:HB1	1:B:282:PRO:HD2	1.69	0.72
1:B:277:TRP:CD1	1:B:277:TRP:N	2.56	0.72
1:A:252:VAL:HB	1:A:279:VAL:CG1	2.22	0.70
1:A:388:SER:O	1:A:389:LYS:HB2	1.90	0.70
1:A:331:ALA:HB1	1:A:336:GLU:CD	2.18	0.69
1:B:189:ARG:HD3	1:B:195:ARG:HD2	1.75	0.68
1:A:375:ARG:H	1:A:375:ARG:CD	1.98	0.68
1:B:195:ARG:HG2	1:B:195:ARG:NH1	2.07	0.67
1:A:296:LEU:HB3	1:A:301:ILE:CD1	2.24	0.67
1:B:290:CYS:HB2	1:B:291:GLY:HA2	1.76	0.67
1:B:181:GLU:HG3	1:B:185:GLN:HE21	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:GLU:HG3	1:A:304:ARG:NE	2.10	0.67
1:B:357:ARG:O	1:B:361:SER:HB2	1.94	0.67
1:B:296:LEU:HB3	1:B:301:ILE:CD1	2.25	0.66
1:A:291:GLY:C	1:A:293:LEU:H	2.05	0.65
2:A:401:OK6:H3	2:A:401:OK6:N7	2.11	0.64
1:A:131:GLU:O	1:A:132:ASP:HB2	1.96	0.64
1:B:331:ALA:HB1	1:B:336:GLU:CD	2.22	0.64
1:B:232:ARG:HG3	1:B:232:ARG:NH1	2.12	0.63
1:A:255:ARG:HH21	1:A:277:TRP:HB2	1.61	0.63
1:A:357:ARG:O	1:A:361:SER:HB2	1.98	0.63
1:A:190:HIS:CD2	1:A:192:ASN:H	2.04	0.63
1:A:373:MET:HB3	1:A:375:ARG:HH21	1.63	0.62
1:B:359:LEU:HD22	1:B:363:LEU:HD22	1.81	0.62
1:A:336:GLU:CD	1:B:336:GLU:CD	2.67	0.62
1:A:180:ARG:C	1:A:183:GLU:OE1	2.43	0.62
1:B:274:ASP:O	1:B:275:PHE:HB2	1.99	0.61
1:A:235:THR:O	1:A:239:GLU:HG3	2.00	0.61
1:A:359:LEU:HD22	1:A:363:LEU:HD22	1.81	0.61
1:B:190:HIS:HD2	1:B:192:ASN:H	1.47	0.61
1:B:292:THR:OG1	1:B:293:LEU:N	2.33	0.61
1:A:140:GLY:HA3	2:A:401:OK6:C16	2.31	0.60
1:B:331:ALA:HB1	1:B:336:GLU:OE1	2.01	0.59
1:A:142:GLY:HA2	1:A:145:GLY:O	2.02	0.59
1:B:180:ARG:NH2	1:B:183:GLU:HG2	2.17	0.59
1:A:180:ARG:O	1:A:183:GLU:OE1	2.21	0.58
1:A:200:PHE:CE2	1:A:207:TYR:CD2	2.92	0.58
1:B:143:LYS:CE	1:B:143:LYS:H	2.17	0.58
1:B:367:ASN:HD22	1:B:370:GLN:HG3	1.70	0.57
1:A:179:ARG:C	1:A:183:GLU:OE2	2.47	0.57
1:B:232:ARG:HH11	1:B:232:ARG:CG	2.18	0.57
1:B:163:VAL:HG22	1:B:207:TYR:HD1	1.70	0.57
1:A:131:GLU:O	1:A:132:ASP:CB	2.53	0.56
1:A:163:VAL:HG22	1:A:207:TYR:HD1	1.71	0.56
1:A:296:LEU:HB3	1:A:301:ILE:HD13	1.88	0.56
1:A:302:GLU:OE1	1:A:304:ARG:NH2	2.29	0.56
1:B:222:LEU:O	1:B:226:SER:N	2.39	0.55
1:B:200:PHE:CE2	1:B:207:TYR:CD2	2.95	0.55
1:A:317:VAL:HG13	1:A:328:PRO:HD2	1.89	0.55
2:B:401:OK6:H15	2:B:401:OK6:N5	2.22	0.55
1:A:259:PRO:HD3	1:A:295:TYR:CZ	2.41	0.55
1:B:358:ASP:OD2	1:B:380:HIS:HE1	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:GLY:HA3	2:B:401:0K6:C16	2.32	0.54
1:B:317:VAL:HG13	1:B:328:PRO:HD2	1.90	0.54
1:A:166:LYS:O	1:A:167:ALA:C	2.49	0.54
1:B:257:ILE:HG22	1:B:318:LEU:HD11	1.90	0.53
1:A:296:LEU:HB3	1:A:301:ILE:HD11	1.90	0.53
1:A:151:ARG:HD3	1:A:156:LYS:HA	1.90	0.53
1:B:179:ARG:C	1:B:183:GLU:OE2	2.52	0.53
2:A:401:0K6:H15	2:A:401:0K6:N5	2.24	0.53
1:A:166:LYS:HA	1:A:169:LEU:HD12	1.91	0.53
1:A:295:TYR:OH	1:A:321:GLU:OE2	2.25	0.52
1:A:210:LEU:HD13	2:A:401:0K6:H12	1.91	0.52
1:B:380:HIS:O	1:B:381:PRO:C	2.53	0.52
1:B:190:HIS:CD2	1:B:191:PRO:HD2	2.45	0.52
1:B:183:GLU:OE1	1:B:184:ILE:N	2.37	0.52
1:A:302:GLU:HG3	1:A:304:ARG:HE	1.72	0.51
1:A:379:GLU:O	1:A:380:HIS:C	2.54	0.51
1:B:166:LYS:O	1:B:167:ALA:C	2.53	0.51
1:B:143:LYS:HG2	1:B:144:PHE:H	1.76	0.50
1:B:143:LYS:H	1:B:143:LYS:NZ	2.08	0.50
1:A:358:ASP:OD2	1:A:380:HIS:HE1	1.94	0.50
1:B:170:GLU:C	1:B:172:ALA:H	2.18	0.50
1:A:282:PRO:HB2	1:A:305:MET:HE2	1.93	0.50
1:A:336:GLU:CD	1:B:336:GLU:OE1	2.55	0.49
1:B:326:LYS:HD3	1:B:330:GLU:OE1	2.12	0.49
1:A:277:TRP:HE3	1:A:277:TRP:H	1.58	0.49
1:A:302:GLU:HG3	1:A:304:ARG:CZ	2.42	0.49
1:A:251:ARG:NH1	1:A:308:GLU:OE1	2.46	0.49
1:B:180:ARG:CZ	1:B:183:GLU:HG2	2.41	0.49
1:A:301:ILE:HG23	1:A:338:TYR:CD2	2.46	0.49
1:A:147:VAL:HG11	2:A:401:0K6:N7	2.26	0.49
1:A:336:GLU:OE2	1:B:336:GLU:CD	2.55	0.49
1:B:296:LEU:CB	1:B:301:ILE:HD11	2.43	0.49
1:A:296:LEU:HD23	1:A:301:ILE:HD11	1.95	0.49
1:A:247:CYS:O	1:A:252:VAL:HG22	2.12	0.48
1:A:252:VAL:HA	1:A:279:VAL:HB	1.95	0.48
1:B:277:TRP:HA	1:B:278:SER:C	2.36	0.48
1:B:180:ARG:C	1:B:183:GLU:OE1	2.57	0.48
1:B:131:GLU:CD	1:B:131:GLU:H	2.22	0.48
1:B:379:GLU:O	1:B:380:HIS:C	2.57	0.48
1:A:163:VAL:HG22	1:A:207:TYR:CD1	2.49	0.47
1:B:253:ILE:HG21	1:B:308:GLU:HA	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:GLU:CD	1:A:131:GLU:H	2.22	0.47
1:A:180:ARG:HG3	1:A:184:ILE:HD11	1.97	0.47
1:A:184:ILE:O	1:A:188:LEU:HG	2.14	0.47
1:B:147:VAL:HG11	2:B:401:OK6:N7	2.22	0.47
1:A:274:ASP:O	1:A:275:PHE:HB2	2.14	0.47
1:B:200:PHE:HE1	1:B:209:ILE:HD11	1.80	0.47
1:A:259:PRO:HD3	1:A:295:TYR:CE2	2.50	0.47
1:B:255:ARG:HH21	1:B:277:TRP:HB3	1.79	0.47
1:A:192:ASN:HB2	1:A:243:ALA:HB2	1.97	0.46
1:B:292:THR:O	1:B:293:LEU:HB2	2.15	0.46
1:B:284:SER:HB2	1:B:305:MET:HG3	1.97	0.46
1:B:300:MET:HE2	1:B:300:MET:HB3	1.88	0.46
1:A:221:GLU:O	1:A:221:GLU:HG3	2.15	0.46
1:B:143:LYS:H	1:B:143:LYS:HE3	1.79	0.46
2:B:401:OK6:N5	2:B:401:OK6:C9	2.76	0.46
1:A:278:SER:HB3	1:A:279:VAL:H	1.61	0.46
1:A:291:GLY:C	1:A:293:LEU:N	2.69	0.46
1:A:269:GLU:H	1:A:269:GLU:HG2	1.35	0.46
1:A:290:CYS:O	1:A:292:THR:N	2.48	0.45
1:A:307:ASP:HB2	1:A:308:GLU:OE2	2.16	0.45
1:B:337:THR:O	1:B:341:ILE:HG13	2.16	0.45
1:A:257:ILE:HD13	1:A:257:ILE:H	1.81	0.45
1:B:135:ILE:N	1:B:135:ILE:HD13	2.32	0.45
1:A:301:ILE:CG2	1:A:338:TYR:CE2	2.94	0.45
1:A:308:GLU:CD	1:A:308:GLU:H	2.22	0.45
2:A:401:OK6:H2	2:A:401:OK6:H3	1.81	0.45
1:B:269:GLU:H	1:B:269:GLU:HG2	1.56	0.45
1:B:387:SER:OG	1:B:388:SER:N	2.49	0.45
1:B:170:GLU:C	1:B:172:ALA:N	2.75	0.45
1:B:290:CYS:CB	1:B:291:GLY:HA2	2.43	0.45
1:A:143:LYS:HD2	3:A:402:PO4:O4	2.17	0.45
1:A:284:SER:HB3	1:A:285:ARG:H	1.49	0.45
1:A:380:HIS:CD2	1:A:382:TRP:H	2.34	0.44
1:A:222:LEU:O	1:A:226:SER:N	2.51	0.44
1:A:293:LEU:O	1:A:296:LEU:HB2	2.17	0.44
1:B:359:LEU:HD21	1:B:377:VAL:HG22	1.99	0.44
1:A:374:LEU:H	1:A:375:ARG:HH21	1.65	0.44
1:A:253:ILE:HG21	1:A:308:GLU:HA	2.00	0.44
1:A:252:VAL:HB	1:A:279:VAL:CB	2.48	0.44
1:B:315:LEU:HD12	1:B:315:LEU:HA	1.68	0.44
1:B:358:ASP:O	1:B:362:ARG:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:TYR:HH	1:A:321:GLU:CD	2.26	0.43
1:A:366:HIS:O	1:A:368:PRO:HD3	2.18	0.43
1:B:296:LEU:CB	1:B:301:ILE:CD1	2.96	0.43
1:A:374:LEU:H	1:A:375:ARG:NH2	2.16	0.43
1:A:373:MET:HB3	1:A:375:ARG:NH2	2.31	0.43
1:B:131:GLU:O	1:B:132:ASP:CB	2.66	0.43
1:B:170:GLU:O	1:B:172:ALA:N	2.51	0.43
1:B:320:TYR:O	1:B:324:VAL:HG23	2.19	0.43
1:B:255:ARG:NH2	1:B:277:TRP:HB3	2.34	0.43
1:A:241:ALA:O	1:A:245:SER:HB2	2.19	0.43
1:A:265:GLY:HA3	1:A:269:GLU:HG3	2.00	0.43
1:A:298:PRO:O	1:A:302:GLU:HG2	2.19	0.43
1:B:295:TYR:OH	1:B:321:GLU:OE2	2.30	0.43
1:B:235:THR:O	1:B:239:GLU:HG3	2.17	0.43
1:B:253:ILE:O	1:B:278:SER:HA	2.19	0.43
1:B:183:GLU:N	1:B:183:GLU:OE1	2.49	0.42
1:A:324:VAL:HG13	1:A:349:PRO:HG2	2.02	0.42
1:B:179:ARG:O	1:B:183:GLU:OE2	2.36	0.42
1:B:326:LYS:HA	1:B:327:PRO:HD3	1.94	0.42
1:B:163:VAL:HG22	1:B:207:TYR:CD1	2.51	0.42
1:A:140:GLY:HA3	2:A:401:OK6:H7	1.99	0.42
1:A:331:ALA:HB3	1:A:337:THR:OG1	2.20	0.42
1:A:318:LEU:O	1:A:319:CYS:C	2.62	0.42
1:A:327:PRO:HG2	1:A:330:GLU:HB2	2.02	0.42
1:A:278:SER:OG	1:A:279:VAL:HG13	2.20	0.42
1:A:380:HIS:HD2	1:A:382:TRP:H	1.66	0.42
1:B:151:ARG:HA	1:B:157:PHE:O	2.20	0.41
1:A:139:LEU:O	2:A:401:OK6:H16	2.19	0.41
1:A:165:PHE:CE1	1:A:205:ARG:HD3	2.55	0.41
1:A:213:ALA:HA	1:A:214:PRO:HD3	1.88	0.41
1:B:273:ALA:O	1:B:274:ASP:HB3	2.20	0.41
1:A:299:GLU:OE1	1:A:299:GLU:N	2.48	0.41
1:A:180:ARG:HG3	1:A:184:ILE:CD1	2.51	0.41
1:A:253:ILE:CG2	1:A:308:GLU:HA	2.50	0.41
1:B:248:HIS:CE1	1:B:308:GLU:HB2	2.55	0.41
1:B:255:ARG:NE	1:B:279:VAL:HG22	2.36	0.41
1:B:301:ILE:CG2	1:B:338:TYR:CE2	2.98	0.41
1:A:143:LYS:HB3	1:A:143:LYS:HE2	1.86	0.41
1:A:178:LEU:C	1:A:180:ARG:H	2.28	0.41
1:A:282:PRO:CB	1:A:305:MET:HE2	2.50	0.41
1:B:367:ASN:ND2	1:B:370:GLN:CG	2.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:GLU:O	1:B:132:ASP:HB2	2.21	0.40
1:B:139:LEU:HB2	1:B:147:VAL:HG22	2.04	0.40
1:B:222:LEU:O	1:B:226:SER:HA	2.22	0.40
1:B:281:ALA:CB	1:B:282:PRO:CD	2.82	0.40
1:A:380:HIS:O	1:A:381:PRO:C	2.65	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	254/271 (94%)	214 (84%)	25 (10%)	15 (6%)	1	3
1	B	254/271 (94%)	215 (85%)	27 (11%)	12 (5%)	2	6
All	All	508/542 (94%)	429 (84%)	52 (10%)	27 (5%)	1	5

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	ALA
1	A	284	SER
1	A	292	THR
1	B	278	SER
1	B	284	SER
1	A	132	ASP
1	A	291	GLY
1	B	132	ASP
1	B	276	GLY
1	B	281	ALA
1	A	142	GLY
1	A	276	GLY
1	A	294	ASP

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Mol	Chain	Res	Type
1	B	226	SER
1	B	293	LEU
1	A	143	LYS
1	A	278	SER
1	A	280	HIS
1	B	274	ASP
1	B	292	THR
1	A	274	ASP
1	A	275	PHE
1	A	380	HIS
1	B	143	LYS
1	B	171	LYS
1	B	291	GLY
1	A	279	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	226/238 (95%)	186 (82%)	40 (18%)	2	6
1	B	226/238 (95%)	197 (87%)	29 (13%)	4	15
All	All	452/476 (95%)	383 (85%)	69 (15%)	3	10

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	GLU
1	A	141	LYS
1	A	143	LYS
1	A	147	VAL
1	A	151	ARG
1	A	154	GLN
1	A	164	LEU
1	A	175	GLU
1	A	177	GLN

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Mol	Chain	Res	Type
1	A	178	LEU
1	A	179	ARG
1	A	180	ARG
1	A	183	GLU
1	A	195	ARG
1	A	204	THR
1	A	215	LEU
1	A	220	ARG
1	A	221	GLU
1	A	232	ARG
1	A	245	SER
1	A	247	CYS
1	A	252	VAL
1	A	257	ILE
1	A	277	TRP
1	A	278	SER
1	A	283	SER
1	A	284	SER
1	A	285	ARG
1	A	292	THR
1	A	296	LEU
1	A	301	ILE
1	A	302	GLU
1	A	308	GLU
1	A	344	VAL
1	A	354	GLU
1	A	359	LEU
1	A	361	SER
1	A	363	LEU
1	A	375	ARG
1	A	379	GLU
1	B	141	LYS
1	B	143	LYS
1	B	147	VAL
1	B	156	LYS
1	B	175	GLU
1	B	178	LEU
1	B	183	GLU
1	B	195	ARG
1	B	201	HIS
1	B	204	THR
1	B	221	GLU

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Mol	Chain	Res	Type
1	B	232	ARG
1	B	239	GLU
1	B	245	SER
1	B	252	VAL
1	B	253	ILE
1	B	269	GLU
1	B	277	TRP
1	B	290	CYS
1	B	293	LEU
1	B	296	LEU
1	B	300	MET
1	B	335	GLN
1	B	342	SER
1	B	359	LEU
1	B	361	SER
1	B	363	LEU
1	B	367	ASN
1	B	377	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	ASN
1	A	177	GLN
1	A	185	GLN
1	A	190	HIS
1	A	192	ASN
1	A	201	HIS
1	A	335	GLN
1	A	367	ASN
1	A	380	HIS
1	B	177	GLN
1	B	185	GLN
1	B	190	HIS
1	B	335	GLN
1	B	367	ASN
1	B	370	GLN
1	B	380	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 6 ligands modelled in this entry, 2 are 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$
2	OK6	A	401	-	36,37,37	1.07	3 (8%)	39,54,54	3.04	17 (43%)
2	OK6	B	401	-	36,37,37	0.98	1 (2%)	39,54,54	2.97	18 (46%)
3	PO4	B	402	4	4,4,4	0.66	0	6,6,6	0.89	0
3	PO4	A	402	4	4,4,4	0.65	0	6,6,6	0.61	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
2	OK6	A	401	-	-	5/20/37/37	0/5/5/5
2	OK6	B	401	-	-	4/20/37/37	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	OK6	C12-N6	-2.66	1.34	1.38
2	A	401	OK6	C12-N6	-2.57	1.34	1.38
2	A	401	OK6	C7-N3	-2.04	1.36	1.38
2	A	401	OK6	C18-N6	-2.00	1.34	1.39

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	OK6	C4-N2-C3	7.66	125.22	110.63
2	B	401	OK6	C4-N2-C3	6.88	123.75	110.63
2	A	401	OK6	C7-C1-N1	6.79	128.61	118.45
2	A	401	OK6	N8-C11-N5	-6.22	120.41	126.42
2	B	401	OK6	C7-C1-N1	6.13	127.62	118.45
2	B	401	OK6	C4-N2-C5	5.92	121.91	110.63
2	B	401	OK6	N8-C11-N5	-5.65	120.95	126.42
2	B	401	OK6	C21-C12-N5	-5.52	119.58	126.87
2	A	401	OK6	C21-C12-N5	-5.46	119.66	126.87
2	A	401	OK6	C4-N2-C5	5.16	120.47	110.63
2	A	401	OK6	C22-N8-C11	5.13	122.84	115.81
2	B	401	OK6	C2-C3-N2	-4.69	103.32	110.86
2	A	401	OK6	C2-C3-N2	-4.62	103.43	110.86
2	B	401	OK6	C22-N8-C11	4.51	122.00	115.81
2	B	401	OK6	C12-C21-C20	4.47	111.92	106.86
2	B	401	OK6	C21-C20-C18	-4.44	100.67	107.31
2	A	401	OK6	C2-N1-C6	3.46	119.74	112.68
2	A	401	OK6	C21-C20-C18	-3.46	102.14	107.31
2	A	401	OK6	C12-C21-C20	3.45	110.78	106.86
2	A	401	OK6	C6-C5-N2	-3.36	105.46	110.86
2	B	401	OK6	C3-N2-C5	3.00	114.32	109.54
2	A	401	OK6	C3-N2-C5	2.97	114.28	109.54
2	B	401	OK6	C6-C5-N2	-2.93	106.16	110.86
2	A	401	OK6	C21-C12-N6	-2.92	106.26	109.52
2	B	401	OK6	C21-C12-N6	-2.85	106.34	109.52
2	A	401	OK6	O1-C1-C7	-2.83	111.63	120.36
2	B	401	OK6	C20-C18-N6	2.80	112.20	108.72
2	B	401	OK6	C2-N1-C6	2.76	118.31	112.68
2	B	401	OK6	O1-C1-C7	-2.53	112.56	120.36
2	A	401	OK6	C22-C21-C12	2.50	121.58	117.20
2	A	401	OK6	C13-N6-C18	-2.42	121.28	126.07
2	A	401	OK6	C3-C2-N1	2.25	114.89	110.42
2	B	401	OK6	C20-C18-C19	-2.23	123.87	129.51
2	B	401	OK6	C13-N6-C18	-2.14	121.84	126.07
2	B	401	OK6	C22-C21-C12	2.03	120.76	117.20

There are no chirality outliers.

All (9) torsion outliers are listed below:

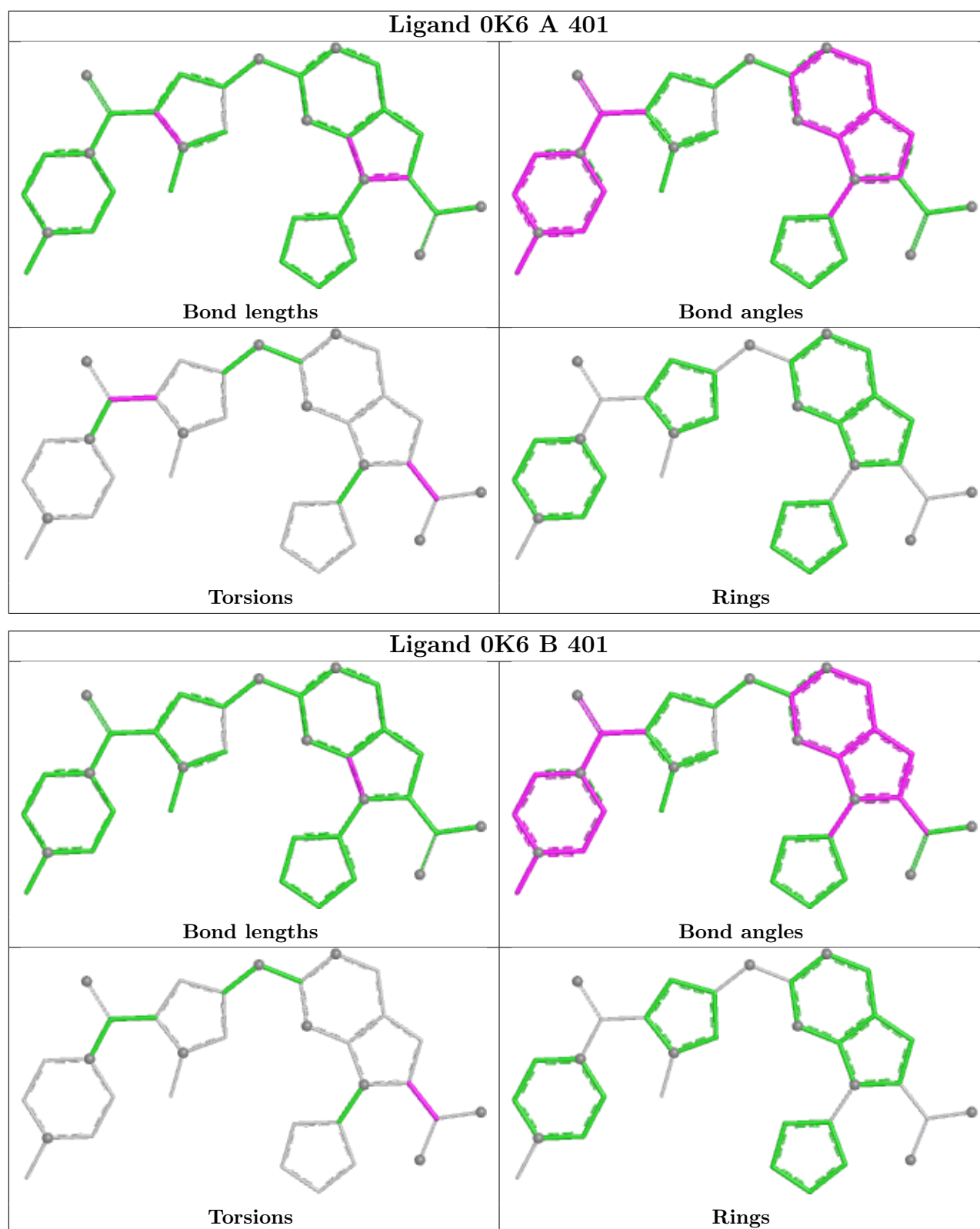
Mol	Chain	Res	Type	Atoms
2	A	401	OK6	C20-C18-C19-O2
2	A	401	OK6	N6-C18-C19-N7
2	A	401	OK6	C20-C18-C19-N7
2	B	401	OK6	N6-C18-C19-O2
2	B	401	OK6	C20-C18-C19-O2
2	B	401	OK6	N6-C18-C19-N7
2	B	401	OK6	C20-C18-C19-N7
2	A	401	OK6	N6-C18-C19-O2
2	A	401	OK6	N1-C1-C7-C23

There are no ring outliers.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	OK6	9	0
2	B	401	OK6	6	0
3	A	402	PO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	258/271 (95%)	-1.34	0 100 100	33, 56, 91, 114	0
1	B	258/271 (95%)	-1.33	0 100 100	33, 56, 88, 106	0
All	All	516/542 (95%)	-1.33	0 100 100	33, 56, 90, 114	0

There are no RSRZ outliers to report.

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](#)

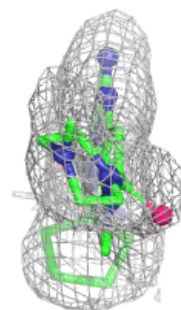
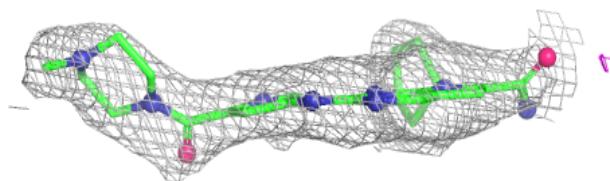
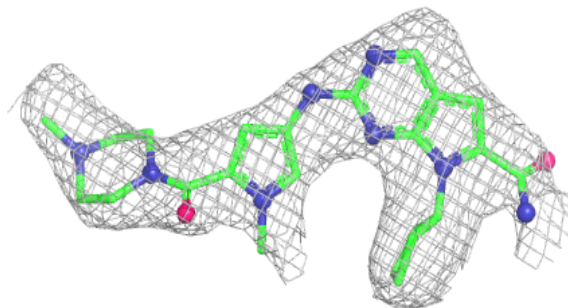
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PO4	A	402	5/5	0.97	0.05	114,120,120,120	0
4	MG	A	403	1/1	0.98	0.05	79,79,79,79	0
2	0K6	A	401	33/33	0.99	0.04	41,57,73,83	0
3	PO4	B	402	5/5	0.99	0.04	93,100,116,117	0
2	0K6	B	401	33/33	0.99	0.04	38,60,74,82	0
4	MG	B	403	1/1	0.99	0.03	68,68,68,68	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.

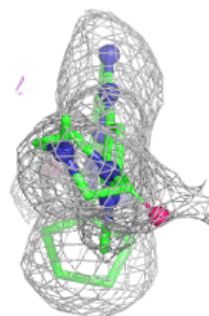
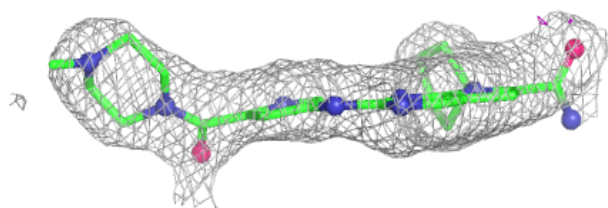
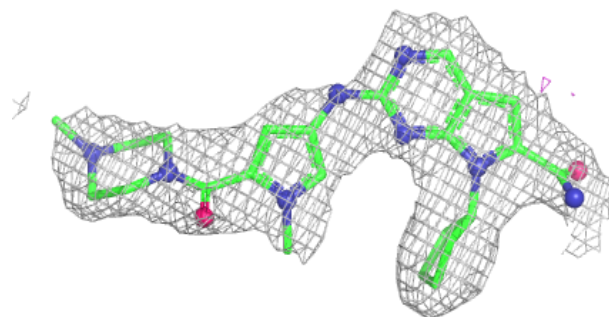
Electron density around 0K6 A 401:

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



Electron density around 0K6 B 401:

$2mF_o-DF_c$ (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.