



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:39 AM UTC

PDB ID : 5OT5 / pdb_00005ot5
Title : The crystal structure of CK2alpha in complex with compound 24
Authors : Brear, P.; De Fusco, C.; Iegre, J.; Yoshida, M.; Mitchell, S.; Rossmann, M.; Carro, L.; Sore, H.; Hyvonen, M.; Spring, D.
Deposited on : 2017-08-21
Resolution : 1.63 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

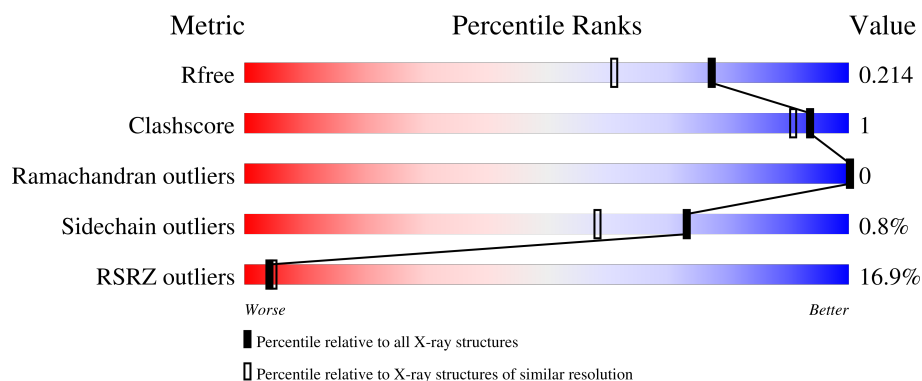
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

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	352	6% 88% 5% 7%
1	B	352	25% 88% • 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6035 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 Casein kinase II subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	10	0
			2833	1812	501	509	11			
1	B	325	Total	C	N	O	S	0	3	0
			2765	1771	483	500	11			

There are 50 discrepancies between the modelled and reference sequences:

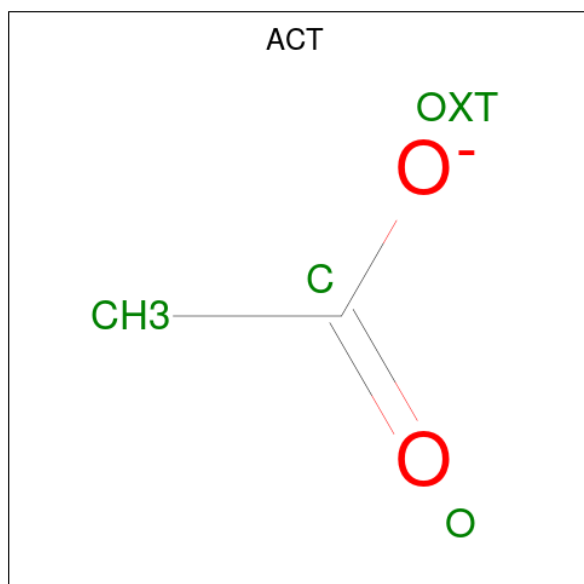
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	GLY	-	expression tag	UNP P68400
A	-21	SER	-	expression tag	UNP P68400
A	-20	MET	-	expression tag	UNP P68400
A	-19	ASP	-	expression tag	UNP P68400
A	-18	ILE	-	expression tag	UNP P68400
A	-17	GLU	-	expression tag	UNP P68400
A	-16	PHE	-	expression tag	UNP P68400
A	-15	ASP	-	expression tag	UNP P68400
A	-14	ASP	-	expression tag	UNP P68400
A	-13	ASP	-	expression tag	UNP P68400
A	-12	ALA	-	expression tag	UNP P68400
A	-11	ASP	-	expression tag	UNP P68400
A	-10	ASP	-	expression tag	UNP P68400
A	-9	ASP	-	expression tag	UNP P68400
A	-8	GLY	-	expression tag	UNP P68400
A	-7	SER	-	expression tag	UNP P68400
A	-6	GLY	-	expression tag	UNP P68400
A	-5	SER	-	expression tag	UNP P68400
A	-4	GLY	-	expression tag	UNP P68400
A	-3	SER	-	expression tag	UNP P68400
A	-2	GLY	-	expression tag	UNP P68400
A	-1	SER	-	expression tag	UNP P68400
A	0	GLY	-	expression tag	UNP P68400
A	1	SER	-	expression tag	UNP P68400
A	21	SER	ARG	engineered mutation	UNP P68400

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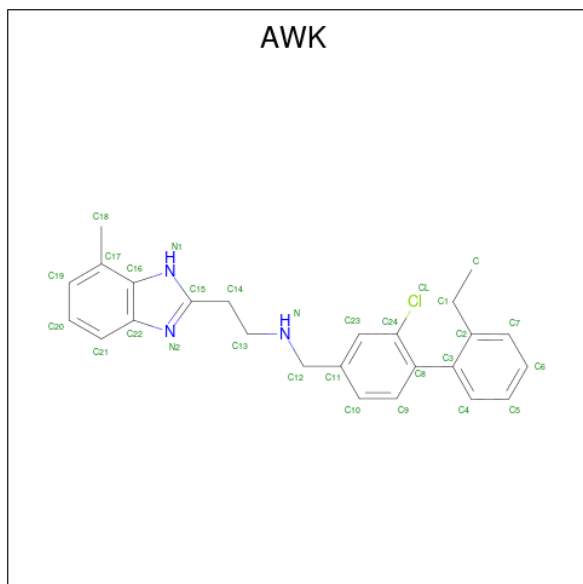
Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	GLY	-	expression tag	UNP P68400
B	-21	SER	-	expression tag	UNP P68400
B	-20	MET	-	expression tag	UNP P68400
B	-19	ASP	-	expression tag	UNP P68400
B	-18	ILE	-	expression tag	UNP P68400
B	-17	GLU	-	expression tag	UNP P68400
B	-16	PHE	-	expression tag	UNP P68400
B	-15	ASP	-	expression tag	UNP P68400
B	-14	ASP	-	expression tag	UNP P68400
B	-13	ASP	-	expression tag	UNP P68400
B	-12	ALA	-	expression tag	UNP P68400
B	-11	ASP	-	expression tag	UNP P68400
B	-10	ASP	-	expression tag	UNP P68400
B	-9	ASP	-	expression tag	UNP P68400
B	-8	GLY	-	expression tag	UNP P68400
B	-7	SER	-	expression tag	UNP P68400
B	-6	GLY	-	expression tag	UNP P68400
B	-5	SER	-	expression tag	UNP P68400
B	-4	GLY	-	expression tag	UNP P68400
B	-3	SER	-	expression tag	UNP P68400
B	-2	GLY	-	expression tag	UNP P68400
B	-1	SER	-	expression tag	UNP P68400
B	0	GLY	-	expression tag	UNP P68400
B	1	SER	-	expression tag	UNP P68400
B	21	SER	ARG	engineered mutation	UNP P68400

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 3 is {N}-[[3-chloranyl-4-(2-ethylphenyl)phenyl]methyl]-2-(7-methyl-1 {H}-benzimidazol-2-yl)ethanamine (CCD ID: AWK) (formula: C₂₅H₂₆ClN₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	0	1
			58	50	2	6		
3	B	1	Total	C	Cl	N	0	0
			29	25	1	3		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	219	Total 219	O 219	0	0
6	B	90	Total 90	O 90	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	64.89Å 68.19Å 332.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.86 – 1.63 23.86 – 1.63	Depositor EDS
% Data completeness (in resolution range)	98.4 (23.86-1.63) 98.3 (23.86-1.63)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 1.63Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.215 , 0.239 (Not available) , 0.214	Depositor DCC
R_{free} test set	4613 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.597	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.077 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6035	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.54% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, PO4, PEG, AWK

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	0.91	1/2909 (0.0%)	1.04	5/3933 (0.1%)
1	B	0.80	1/2840 (0.0%)	1.06	2/3842 (0.1%)
All	All	0.85	2/5749 (0.0%)	1.05	7/7775 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	MET	SD-CE	-7.05	1.61	1.79
1	B	150	MET	SD-CE	-6.19	1.64	1.79

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	ASP	CA-C-N	6.67	129.52	120.38
1	B	120	ASP	C-N-CA	6.67	129.52	120.38
1	A	124	LEU	CA-C-N	6.36	129.44	120.28
1	A	124	LEU	C-N-CA	6.36	129.44	120.28
1	A	55	GLU	N-CA-C	-5.83	100.16	109.96
1	A	197	PHE	CA-CB-CG	-5.14	108.66	113.80
1	A	118	ASN	N-CA-C	5.12	117.98	110.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2833	0	2771	11	0
1	B	2765	0	2696	5	0
2	A	16	0	12	0	0
2	B	8	0	6	0	0
3	A	58	0	0	3	0
3	B	29	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	7	0	10	2	0
6	A	219	0	0	2	1
6	B	90	0	0	0	0
All	All	6035	0	5495	16	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 1.

All (16) 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:196:TYR:HD2	5:A:407:PEG:H32	1.62	0.62
1:A:196:TYR:CD2	5:A:407:PEG:H32	2.38	0.59
1:A:44:LYS:HD3	1:A:47[B]:ARG:HG2	1.87	0.54
1:A:162:VAL:HG11	3:A:405[A]:AWK:CL	2.46	0.53
1:A:225:MET:HE3	3:A:405[B]:AWK:CL	2.46	0.52
1:B:303:LYS:HB3	1:B:313:LEU:HG	1.95	0.48
1:A:59:ILE:HD11	6:A:531:HOH:O	2.13	0.48
1:B:53:VAL:HG13	1:B:66:VAL:HG13	1.97	0.46
1:A:54:PHE:CE1	1:A:69:ILE:HD12	2.53	0.44
1:A:225:MET:CE	3:A:405[B]:AWK:CL	3.03	0.44
1:A:114:GLU:O	1:A:114:GLU:HG3	2.18	0.41
1:B:244:ARG:NH1	1:B:247:LYS:HE3	2.35	0.41
1:A:236:HIS:HE1	6:A:686:HOH:O	2.03	0.41
1:A:95:ILE:HB	1:A:174:ILE:HG22	2.03	0.41
1:B:5:VAL:HB	1:B:261:TYR:HA	2.01	0.41
1:B:95:ILE:HB	1:B:174:ILE:HG22	2.02	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)
6:A:519:HOH:O	6:A:519:HOH:O[4_597]	0.82	1.38

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	305/352 (87%)	294 (96%)	11 (4%)	0	100	100
1	B	289/352 (82%)	282 (98%)	7 (2%)	0	100	100
All	All	594/704 (84%)	576 (97%)	18 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	303/319 (95%)	302 (100%)	1 (0%)	86	79
1	B	302/319 (95%)	298 (99%)	4 (1%)	61	38
All	All	605/638 (95%)	600 (99%)	5 (1%)	73	58

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	VAL
1	B	66	VAL
1	B	73	VAL

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Mol	Chain	Res	Type
1	B	124	LEU
1	B	258	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	HIS
1	A	168	HIS
1	A	186	GLN
1	A	236	HIS
1	A	262	ASN
1	A	270	ASN
1	B	168	HIS
1	B	186	GLN
1	B	207	GLN
1	B	234	HIS
1	B	262	ASN
1	B	270	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](#)

12 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	$\# Z > 2$
5	PEG	A	407	-	6,6,6	0.26	0	5,5,5	0.26	0
2	ACT	B	402	-	3,3,3	1.11	0	3,3,3	0.90	0
2	ACT	A	402	-	3,3,3	1.15	0	3,3,3	1.15	0
4	PO4	A	406	-	4,4,4	2.00	2 (50%)	6,6,6	0.96	0
2	ACT	A	403	-	3,3,3	1.10	0	3,3,3	0.87	0
3	AWK	A	405[B]	-	32,32,32	0.59	0	41,44,44	1.34	3 (7%)
3	AWK	A	405[A]	-	32,32,32	0.71	1 (3%)	41,44,44	1.39	2 (4%)
2	ACT	B	401	-	3,3,3	1.47	0	3,3,3	0.90	0
2	ACT	A	404	-	3,3,3	1.28	1 (33%)	3,3,3	1.09	0
2	ACT	A	401	-	3,3,3	1.63	1 (33%)	3,3,3	1.04	0
4	PO4	B	404	-	4,4,4	1.88	1 (25%)	6,6,6	0.80	0
3	AWK	B	403	-	32,32,32	0.58	0	41,44,44	1.29	2 (4%)

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
5	PEG	A	407	-	-	3/4/4/4	-
3	AWK	A	405[B]	-	-	0/13/13/13	0/4/4/4
3	AWK	B	403	-	-	0/13/13/13	0/4/4/4
3	AWK	A	405[A]	-	-	0/13/13/13	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	405[A]	AWK	C16-C22	-2.78	1.38	1.40
2	A	401	ACT	CH3-C	2.28	1.58	1.49
4	A	406	PO4	P-O2	2.18	1.61	1.54
4	B	404	PO4	P-O3	2.08	1.60	1.54
2	A	404	ACT	O-C	2.06	1.31	1.22
4	A	406	PO4	P-O4	2.05	1.60	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	405[A]	AWK	C16-C22-N2	-6.09	105.75	109.91
3	A	405[B]	AWK	C16-C22-N2	-6.04	105.78	109.91
3	B	403	AWK	C16-C22-N2	-5.94	105.85	109.91
3	A	405[A]	AWK	C22-C16-N1	5.11	109.66	105.40
3	A	405[B]	AWK	C22-C16-N1	4.59	109.22	105.40
3	B	403	AWK	C22-C16-N1	4.28	108.97	105.40
3	A	405[B]	AWK	C14-C15-N2	-2.03	122.64	125.03

There are no chirality outliers.

All (3) torsion outliers are listed below:

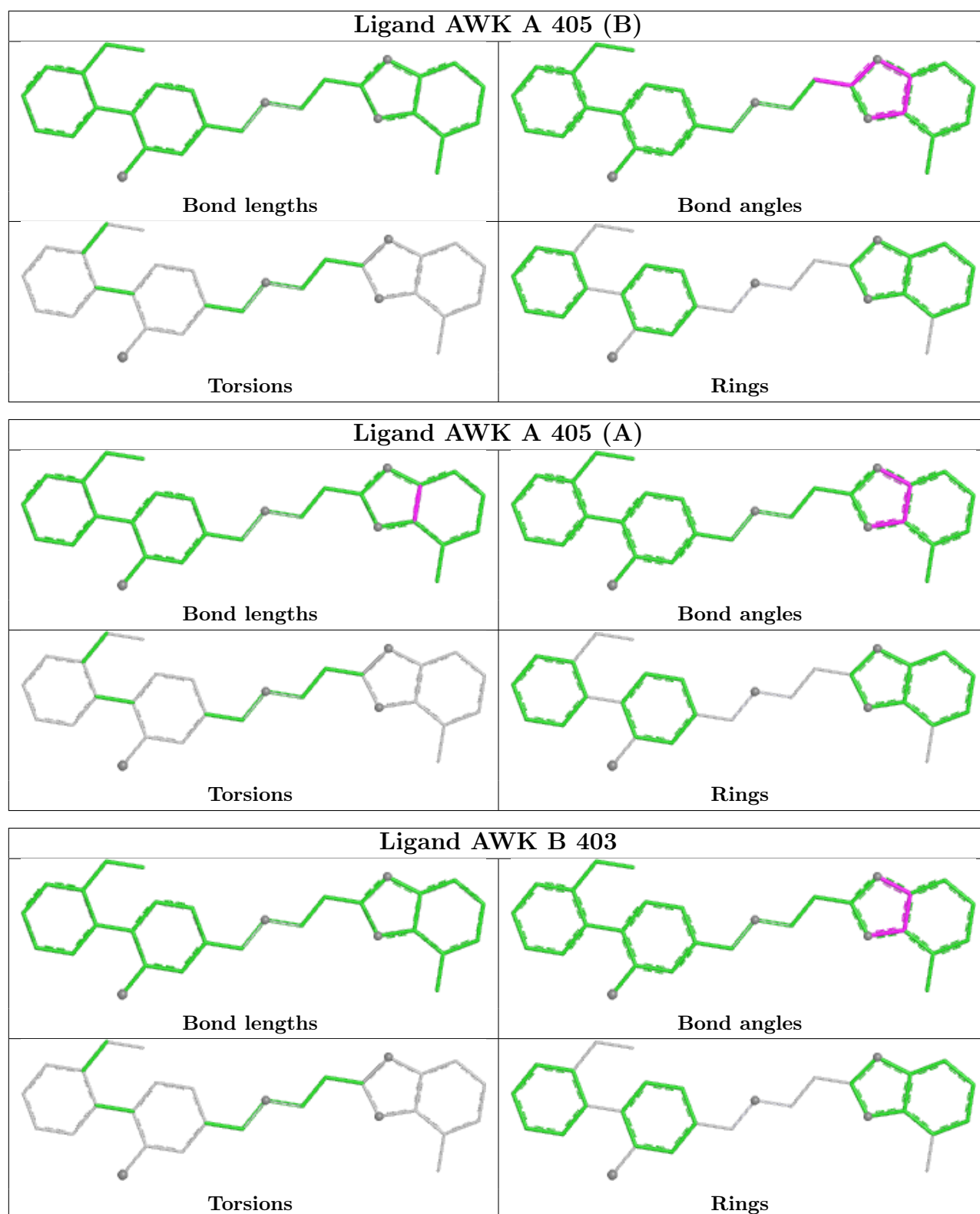
Mol	Chain	Res	Type	Atoms
5	A	407	PEG	C4-C3-O2-C2
5	A	407	PEG	O1-C1-C2-O2
5	A	407	PEG	C1-C2-O2-C3

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	407	PEG	2	0
3	A	405[B]	AWK	2	0
3	A	405[A]	AWK	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	326/352 (92%)	0.35	22 (6%) 24 27	11, 29, 56, 96	10 (3%)
1	B	325/352 (92%)	1.47	88 (27%) 1 1	18, 48, 84, 116	3 (0%)
All	All	651/704 (92%)	0.91	110 (16%) 4 5	11, 39, 79, 116	13 (1%)

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	124	LEU	5.0
1	A	48[A]	GLY	5.0
1	A	50[A]	TYR	5.0
1	A	49[A]	LYS	4.9
1	A	125	TYR	4.9
1	B	327	VAL	4.8
1	B	50	TYR	4.5
1	B	121	PHE	4.5
1	B	307	TYR	4.5
1	B	255	TYR	4.4
1	B	45	LEU	4.4
1	B	263	ILE	4.2
1	B	128	LEU	4.0
1	B	5	VAL	4.0
1	A	328	VAL	4.0
1	B	73	VAL	4.0
1	B	257	TYR	3.9
1	A	73	VAL	3.9
1	B	53	VAL	3.8
1	B	248	VAL	3.8
1	A	124	LEU	3.7
1	B	249	LEU	3.7
1	B	273	LEU	3.6
1	A	33	TRP	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	46	GLY	3.6
1	B	282	GLU	3.6
1	B	251	THR	3.5
1	B	125	TYR	3.5
1	B	272	ILE	3.5
1	A	105	VAL	3.4
1	B	269	PHE	3.4
1	B	258	ILE	3.3
1	B	323	TYR	3.3
1	B	42	VAL	3.3
1	B	72	PRO	3.2
1	B	129	THR	3.2
1	B	297	ALA	3.1
1	B	284	PHE	3.1
1	B	281	TRP	3.1
1	B	206	TYR	3.1
1	B	265	LEU	3.0
1	B	48	GLY	3.0
1	A	74	LYS	3.0
1	B	233	PHE	2.9
1	B	4	PRO	2.9
1	B	123	GLN	2.9
1	B	204	VAL	2.8
1	B	245	ILE	2.8
1	A	121	PHE	2.8
1	B	216	TRP	2.7
1	B	227	PHE	2.7
1	B	292	LEU	2.7
1	B	59	ILE	2.7
1	B	41	LEU	2.7
1	B	30	VAL	2.6
1	A	107[A]	ARG	2.6
1	A	123	GLN	2.5
1	B	22[A]	GLU	2.5
1	B	31	VAL	2.5
1	B	293	VAL	2.5
1	B	276	HIS	2.5
1	B	300	PHE	2.5
1	B	122	LYS	2.4
1	B	298	LEU	2.4
1	B	325	TYR	2.4
1	B	285	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	120	ASP	2.4
1	B	3	GLY	2.4
1	A	72	PRO	2.4
1	B	226	ILE	2.4
1	A	31	VAL	2.4
1	B	43	ARG	2.3
1	B	254	LEU	2.3
1	B	18	HIS	2.3
1	B	54	PHE	2.3
1	A	104	PRO	2.3
1	B	225	MET	2.3
1	B	203	LEU	2.3
1	B	305	LEU	2.3
1	A	22	GLU	2.3
1	B	261	TYR	2.3
1	B	33	TRP	2.2
1	B	267	PRO	2.2
1	B	301	LEU	2.2
1	B	205	ASP	2.2
1	B	283	ARG	2.2
1	B	75	LYS	2.2
1	A	59	ILE	2.2
1	B	78	ILE	2.2
1	B	105	VAL	2.2
1	B	318	ALA	2.2
1	A	128	LEU	2.2
1	A	47[A]	ARG	2.2
1	B	69	ILE	2.2
1	B	17	THR	2.2
1	B	271	ASP	2.1
1	A	327	VAL	2.1
1	B	287	SER	2.1
1	B	119	THR	2.1
1	B	82	ILE	2.1
1	B	242	LEU	2.1
1	B	326	THR	2.1
1	B	291	HIS	2.1
1	B	246	ALA	2.1
1	B	256	ASP	2.1
1	B	127	THR	2.1
1	B	243	VAL	2.0
1	B	280	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	230	GLU	2.0
1	B	274	GLY	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](#)

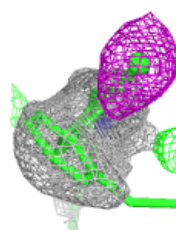
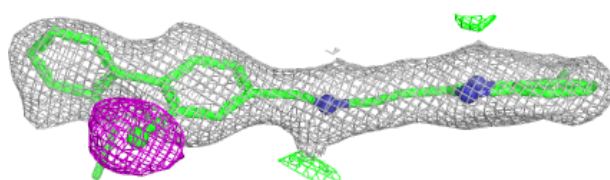
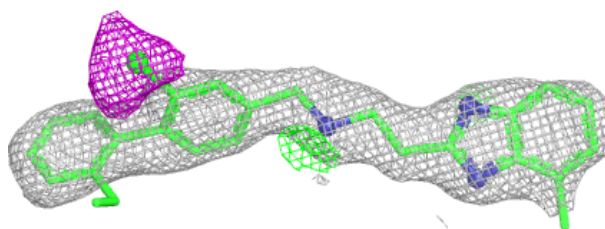
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	ACT	A	404	4/4	0.75	0.20	44,45,46,51	0
3	AWK	B	403	29/29	0.75	0.15	52,58,72,80	0
5	PEG	A	407	7/7	0.82	0.12	32,40,49,52	0
3	AWK	A	405[A]	29/29	0.84	0.12	22,35,47,48	29
3	AWK	A	405[B]	29/29	0.84	0.12	32,42,51,56	29
4	PO4	B	404	5/5	0.88	0.09	48,51,52,54	0
2	ACT	A	403	4/4	0.88	0.12	38,39,41,44	0
2	ACT	B	401	4/4	0.92	0.10	35,35,39,39	0
4	PO4	A	406	5/5	0.92	0.08	38,39,41,41	0
2	ACT	B	402	4/4	0.93	0.09	52,53,53,53	0
2	ACT	A	401	4/4	0.96	0.06	28,29,30,30	0
2	ACT	A	402	4/4	0.96	0.09	30,36,38,39	0

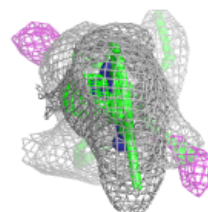
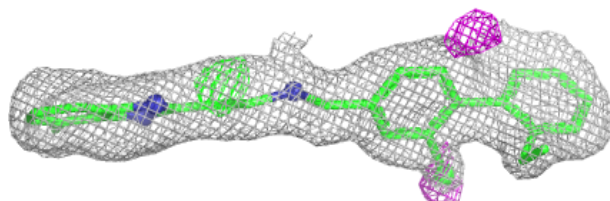
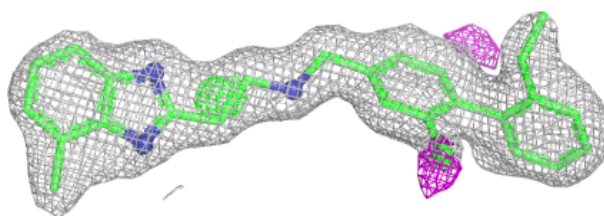
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AWK B 403:

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

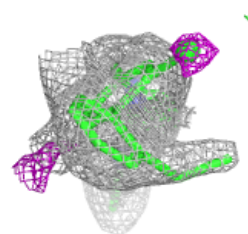
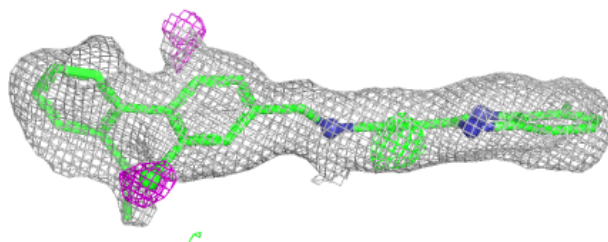
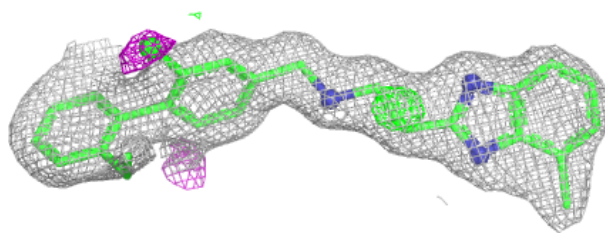
**Electron density around AWK A 405 (A):**

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



Electron density around AWK A 405 (B):

$2mF_o-DF_c$ (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.