



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2024 – 01:17 PM EDT

PDB ID : 5F9E
Title : Structure of Protein Kinase C theta with compound 10: 2,2-dimethyl-7-(2-oxidanylidene-3 {H}-imidazo[4,5-b]pyridin-1-yl)-1-(phenylmethyl)-3 {H}-quinazolin-4-one
Authors : Klein, M.
Deposited on : 2015-12-09
Resolution : 2.00 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

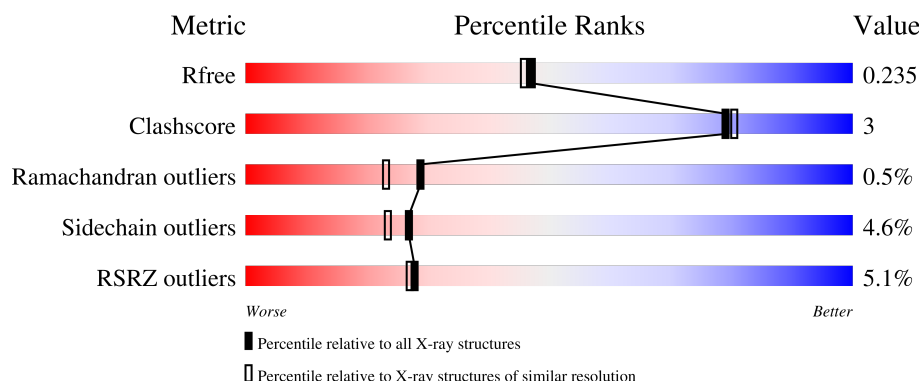
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	353	<div> <div>2%</div> <div>87%</div> <div>7% • 5%</div> </div>
1	B	353	<div> <div>7%</div> <div>73%</div> <div>9% • 17%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5508 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.

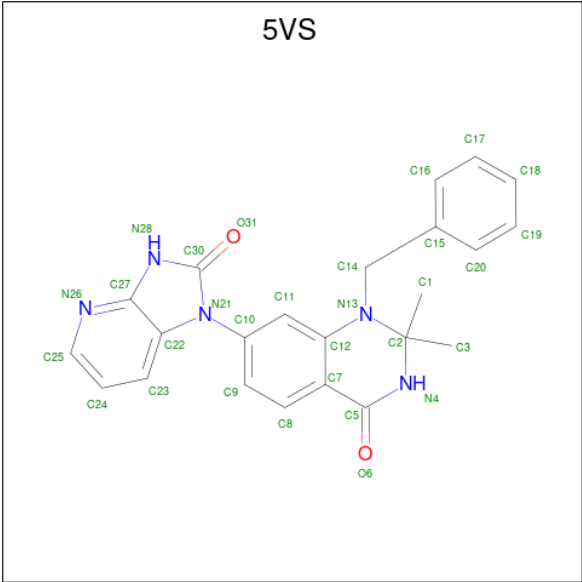
- Molecule 1 is a protein called Protein kinase C theta type.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	P	S	0	6	0
			2832	1831	474	503	2	22			
1	B	294	Total	C	N	O	P	S	0	1	0
			2452	1595	406	433	1	17			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	360	MET	-	initiating methionine	UNP Q04759
A	381	GLU	ILE	conflict	UNP Q04759
A	538	GLU	THR	conflict	UNP Q04759
A	707	HIS	-	expression tag	UNP Q04759
A	708	HIS	-	expression tag	UNP Q04759
A	709	HIS	-	expression tag	UNP Q04759
A	710	HIS	-	expression tag	UNP Q04759
A	711	HIS	-	expression tag	UNP Q04759
A	712	HIS	-	expression tag	UNP Q04759
B	360	MET	-	initiating methionine	UNP Q04759
B	381	GLU	ILE	conflict	UNP Q04759
B	538	GLU	THR	conflict	UNP Q04759
B	707	HIS	-	expression tag	UNP Q04759
B	708	HIS	-	expression tag	UNP Q04759
B	709	HIS	-	expression tag	UNP Q04759
B	710	HIS	-	expression tag	UNP Q04759
B	711	HIS	-	expression tag	UNP Q04759
B	712	HIS	-	expression tag	UNP Q04759

- Molecule 2 is 2,2-dimethyl-7-(2-oxidanylidene-3 {H}-imidazo[4,5-b]pyridin-1-yl)-1-(phenylmethyl)-3 {H}-quinazolin-4-one (three-letter code: 5VS) (formula: C₂₃H₂₁N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			30	23	5	2		
2	B	1	Total	C	N	O	0	0
			30	23	5	2		


- Molecule 3 is water.

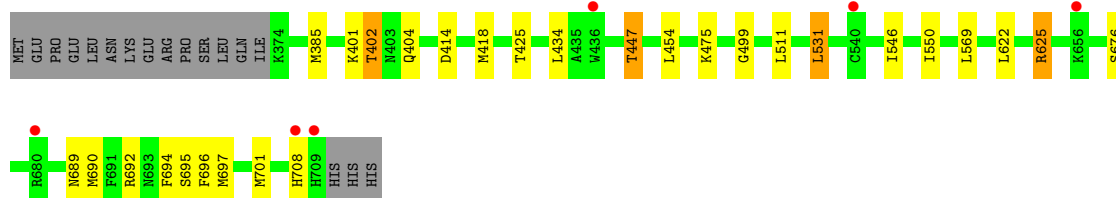
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	126	Total	O	0	0
			126	126		
3	B	38	Total	O	0	0
			38	38		

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.

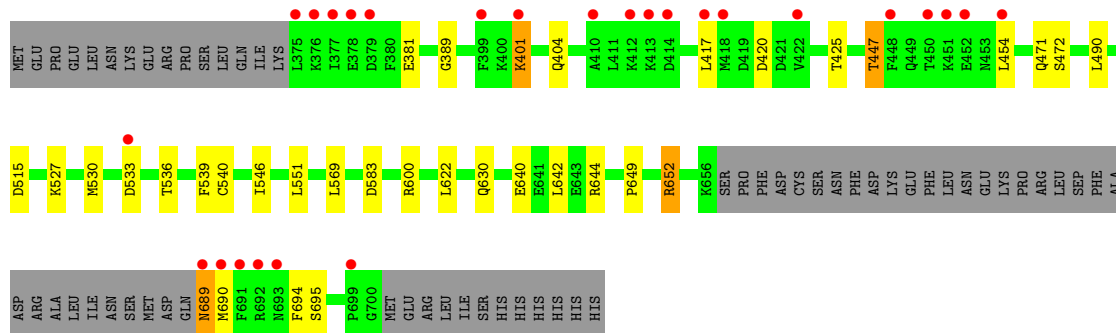
- Molecule 1: Protein kinase C theta type

Chain A: 



- Molecule 1: Protein kinase C theta type

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.60Å 76.79Å 149.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.80 – 2.00 41.76 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.3 (41.80-2.00) 97.4 (41.76-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.197 , 0.233 0.203 , 0.235	Depositor DCC
R_{free} test set	2929 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtriage
Anisotropy	0.491	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5508	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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: 5VS, SEP

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	0.54	0/2903	0.71	0/3898
1	B	0.45	1/2509 (0.0%)	0.65	3/3371 (0.1%)
All	All	0.50	1/5412 (0.0%)	0.69	3/7269 (0.0%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	540	CYS	CB-SG	-5.84	1.72	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	583	ASP	CB-CG-OD1	5.97	123.67	118.30
1	B	540	CYS	N-CA-CB	-5.58	100.55	110.60
1	B	652	ARG	NE-CZ-NH2	-5.20	117.70	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	539	PHE	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	2832	0	2797	19	0
1	B	2452	0	2422	11	0
2	A	30	0	0	0	0
2	B	30	0	0	0	0
3	A	126	0	0	3	0
3	B	38	0	0	1	0
All	All	5508	0	5219	30	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 3.

All (30) 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:625[B]:ARG:HG2	1:A:625[B]:ARG:HH21	1.35	0.92
1:A:625[B]:ARG:HG2	1:A:625[B]:ARG:NH2	1.87	0.87
1:A:625[B]:ARG:HH21	1:A:625[B]:ARG:CG	1.89	0.86
1:A:385:MET:HE3	3:A:1157:HOH:O	1.79	0.81
1:A:418[B]:MET:HE1	3:A:1130:HOH:O	1.80	0.81
1:A:447:THR:CG2	1:A:694:PHE:O	2.47	0.62
1:A:499:GLY:CA	1:A:531:LEU:HD13	2.32	0.59
1:A:546[A]:ILE:HD11	1:A:550:ILE:HG21	1.85	0.58
1:A:499:GLY:HA2	1:A:531:LEU:HD13	1.88	0.56
1:A:414:ASP:O	1:A:418[B]:MET:HE2	2.05	0.55
1:A:447:THR:HG23	1:A:694:PHE:O	2.07	0.55
1:B:689:ASN:HD22	1:B:689:ASN:N	2.08	0.52
1:B:389:GLY:HA3	3:B:1132:HOH:O	2.09	0.52
1:A:418[A]:MET:HA	1:A:418[A]:MET:HE2	1.92	0.52
1:B:640:GLU:O	1:B:644:ARG:HG3	2.10	0.51
1:B:425:THR:HG21	1:B:454:LEU:HD13	1.96	0.48
1:A:447:THR:HG22	1:A:696:PHE:N	2.29	0.47
1:A:531:LEU:HB2	3:A:1199:HOH:O	2.15	0.47
1:B:515:ASP:O	1:B:649:PRO:HG2	2.15	0.46
1:B:447:THR:CG2	1:B:694:PHE:O	2.64	0.45
1:B:527:LYS:HG2	1:B:530:MET:HE3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:697:MET:HE2	1:A:701:MET:HG3	1.98	0.44
1:A:425:THR:HG21	1:A:454:LEU:HD13	2.00	0.43
1:B:515:ASP:OD1	1:B:652:ARG:NH2	2.51	0.43
1:A:546[A]:ILE:HD11	1:A:550:ILE:CG2	2.48	0.43
1:B:490:LEU:HD11	1:B:642:LEU:HD23	2.01	0.42
1:A:402:THR:HG22	1:A:708:HIS:NE2	2.34	0.42
1:A:401:LYS:O	1:A:402:THR:O	2.38	0.41
1:B:546[A]:ILE:HG21	1:B:551:LEU:HD21	2.02	0.41
1:B:530:MET:HE3	1:B:536:THR:HB	2.02	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	338/353 (96%)	330 (98%)	6 (2%)	2 (1%)	25	19
1	B	290/353 (82%)	277 (96%)	12 (4%)	1 (0%)	41	37
All	All	628/706 (89%)	607 (97%)	18 (3%)	3 (0%)	29	23

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	689	ASN
1	A	402	THR
1	B	401	LYS

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	309/320 (97%)	297 (96%)	12 (4%)	32	30
1	B	265/320 (83%)	250 (94%)	15 (6%)	20	16
All	All	574/640 (90%)	547 (95%)	27 (5%)	27	22

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	404	GLN
1	A	434	LEU
1	A	447	THR
1	A	475	LYS
1	A	511	LEU
1	A	531	LEU
1	A	569	LEU
1	A	622	LEU
1	A	625[A]	ARG
1	A	625[B]	ARG
1	A	690	MET
1	A	692	ARG
1	B	381	GLU
1	B	401	LYS
1	B	404	GLN
1	B	417	LEU
1	B	420	ASP
1	B	447	THR
1	B	471	GLN
1	B	472	SER
1	B	533	ASP
1	B	569	LEU
1	B	600	ARG
1	B	622	LEU
1	B	630	GLN
1	B	689	ASN
1	B	690	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	404	GLN

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Mol	Chain	Res	Type
1	A	684	ASN
1	A	693	ASN
1	B	404	GLN
1	B	689	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	SEP	A	695	1	8,9,10	0.96	0	8,12,14	1.72	3 (37%)
1	SEP	B	695	1	8,9,10	0.65	0	8,12,14	1.24	1 (12%)
1	SEP	A	676	1	8,9,10	0.90	0	8,12,14	5.41	2 (25%)

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
1	SEP	A	695	1	-	1/5/8/10	-
1	SEP	B	695	1	-	1/5/8/10	-
1	SEP	A	676	1	-	4/5/8/10	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	676	SEP	OG-CB-CA	14.25	122.01	108.14
1	A	676	SEP	P-OG-CB	-4.79	105.11	118.30
1	A	695	SEP	O3P-P-OG	-2.55	99.94	106.73
1	A	695	SEP	O3P-P-O2P	2.27	116.32	107.64
1	A	695	SEP	O2P-P-OG	-2.26	100.73	106.73
1	B	695	SEP	O3P-P-O2P	2.05	115.49	107.64

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	676	SEP	N-CA-CB-OG
1	A	676	SEP	CB-OG-P-O2P
1	A	676	SEP	CB-OG-P-O3P
1	A	676	SEP	CB-OG-P-O1P
1	A	695	SEP	N-CA-CB-OG
1	B	695	SEP	N-CA-CB-OG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	5VS	B	1001	-	34,34,34	0.98	2 (5%)	42,51,51	2.61	12 (28%)
2	5VS	A	1001	-	34,34,34	1.04	3 (8%)	42,51,51	2.46	11 (26%)

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	5VS	B	1001	-	-	2/8/27/27	0/5/5/5
2	5VS	A	1001	-	-	2/8/27/27	0/5/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	5VS	C27-N28	2.69	1.40	1.37
2	A	1001	5VS	C10-N21	-2.57	1.40	1.44
2	A	1001	5VS	C3-C2	2.38	1.56	1.52
2	B	1001	5VS	C30-N28	-2.27	1.34	1.37
2	B	1001	5VS	C30-N21	-2.24	1.36	1.40

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	5VS	N28-C30-N21	8.93	110.62	106.37
2	A	1001	5VS	N28-C30-N21	8.20	110.28	106.37
2	B	1001	5VS	C3-C2-C1	-7.81	104.85	111.71
2	A	1001	5VS	O6-C5-N4	-7.67	115.71	121.74
2	B	1001	5VS	C22-N21-C30	-5.01	106.67	109.24
2	B	1001	5VS	O31-C30-N28	-4.78	123.23	127.50
2	A	1001	5VS	C15-C14-N13	3.87	118.89	114.16
2	B	1001	5VS	O6-C5-N4	-3.75	118.80	121.74
2	A	1001	5VS	O31-C30-N28	-3.73	124.16	127.50
2	B	1001	5VS	C15-C14-N13	-3.67	109.67	114.16
2	A	1001	5VS	C22-N21-C30	-3.46	107.46	109.24
2	A	1001	5VS	C1-C2-N4	-3.44	102.37	109.88
2	B	1001	5VS	N28-C27-N26	3.43	132.74	125.61
2	A	1001	5VS	N28-C27-N26	3.39	132.66	125.61
2	B	1001	5VS	C12-C7-C5	-3.36	117.50	120.08
2	A	1001	5VS	C7-C5-N4	3.33	120.42	115.91
2	A	1001	5VS	C3-C2-C1	-3.26	108.85	111.71
2	A	1001	5VS	C25-N26-C27	2.37	121.46	116.77
2	B	1001	5VS	C25-N26-C27	2.36	121.46	116.77
2	A	1001	5VS	C12-C7-C5	-2.28	118.32	120.08
2	B	1001	5VS	C7-C5-N4	2.21	118.91	115.91
2	B	1001	5VS	C9-C10-N21	2.07	121.97	119.74
2	B	1001	5VS	C1-C2-N4	-2.07	105.36	109.88

There are no chirality outliers.

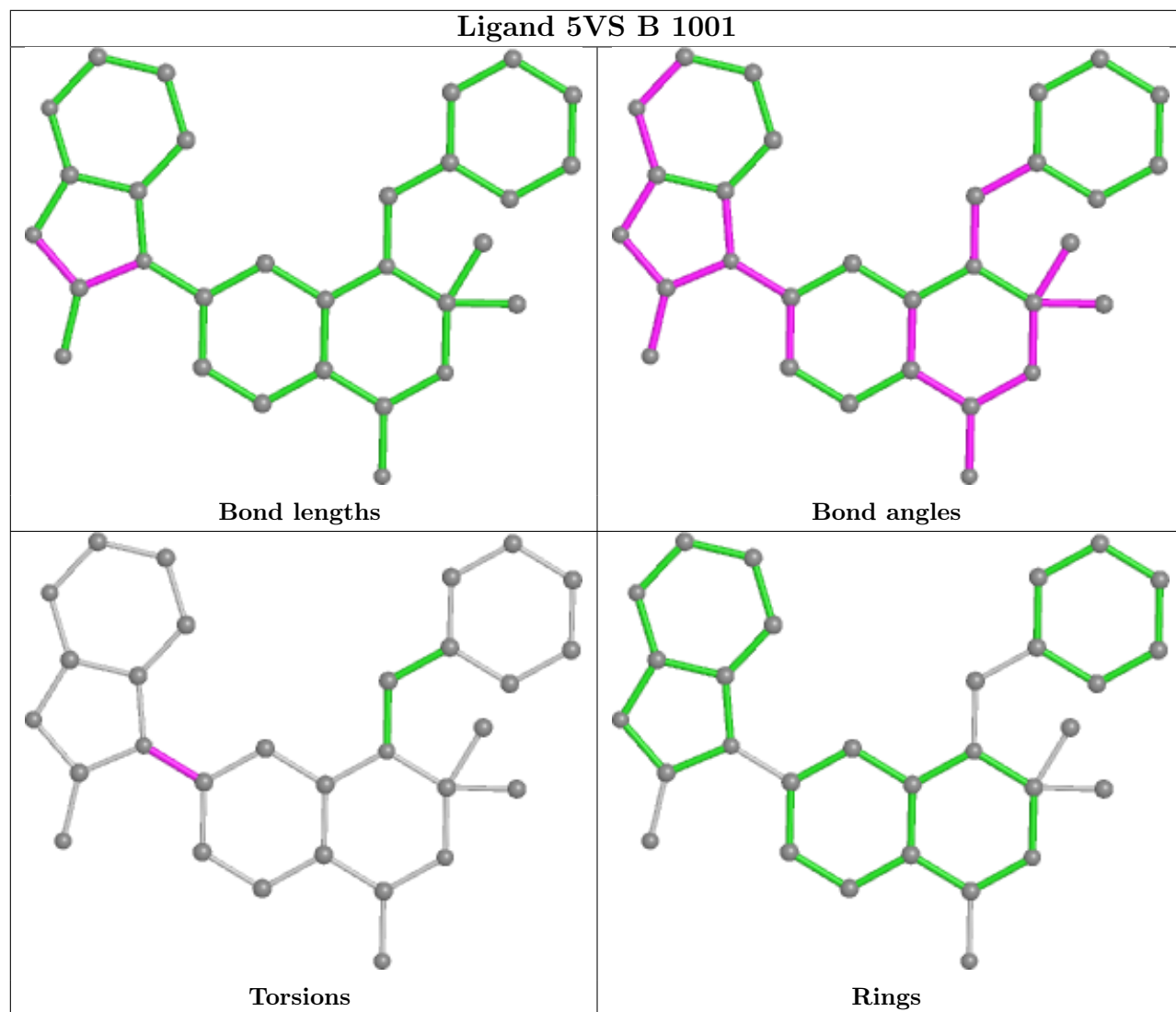
All (4) torsion outliers are listed below:

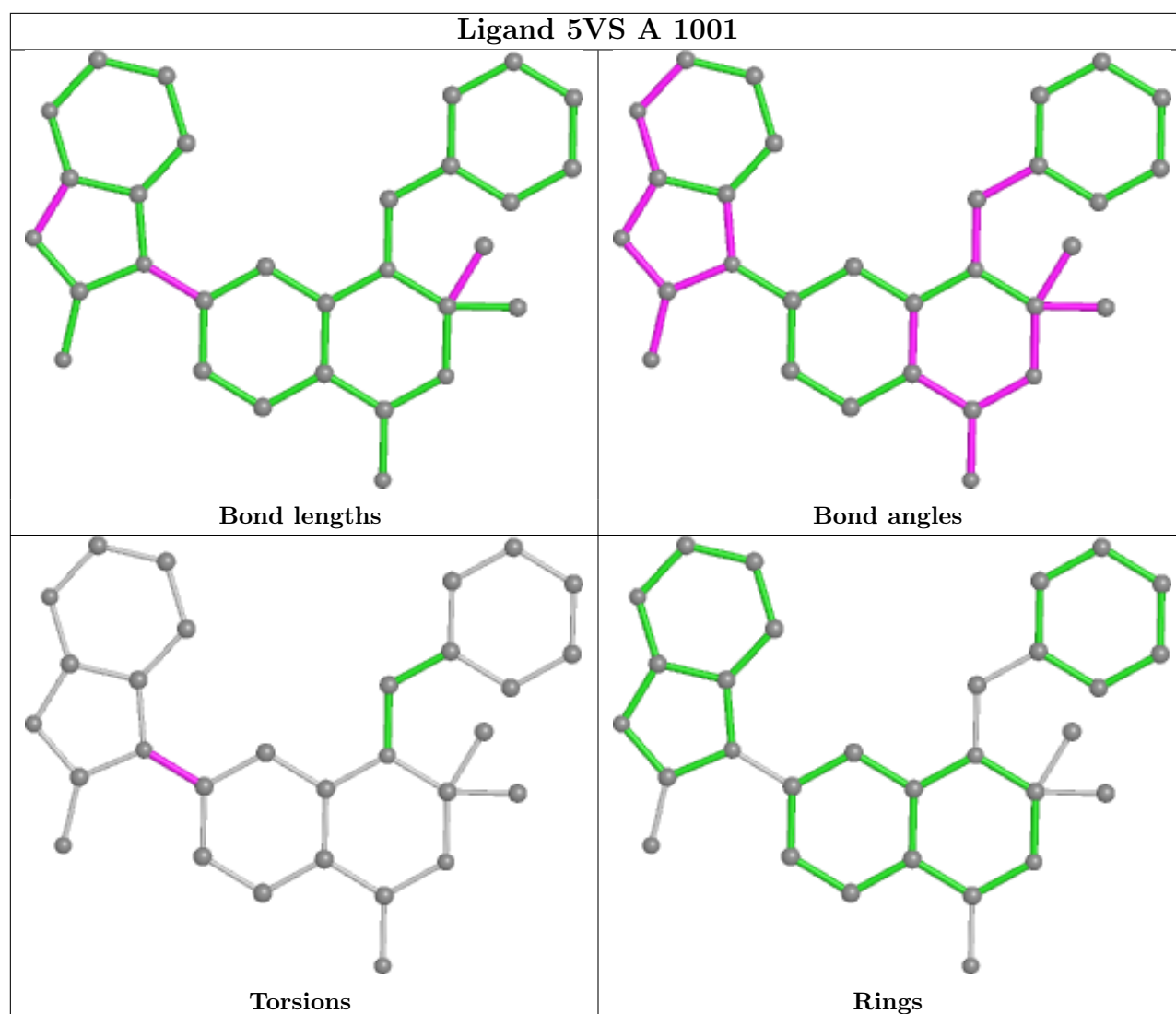
Mol	Chain	Res	Type	Atoms
2	B	1001	5VS	C11-C10-N21-C22
2	B	1001	5VS	C9-C10-N21-C22
2	A	1001	5VS	C11-C10-N21-C22
2	A	1001	5VS	C9-C10-N21-C22

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	334/353 (94%)	0.17	6 (1%) 68 66	29, 43, 72, 101	0
1	B	293/353 (83%)	0.49	26 (8%) 9 8	38, 60, 102, 119	0
All	All	627/706 (88%)	0.32	32 (5%) 28 27	29, 51, 90, 119	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	451	LYS	6.2
1	B	690	MET	5.8
1	B	692	ARG	5.4
1	B	375	LEU	5.4
1	B	454	LEU	5.2
1	B	417	LEU	4.8
1	B	693	ASN	4.7
1	B	376	LYS	4.0
1	B	399	PHE	3.7
1	A	436	TRP	3.7
1	A	709	HIS	3.3
1	B	418	MET	3.3
1	B	401	LYS	3.3
1	B	691	PHE	3.2
1	A	680	ARG	3.1
1	B	450	THR	3.0
1	B	689	ASN	2.9
1	B	377	ILE	2.7
1	B	699	PRO	2.6
1	B	378	GLU	2.6
1	B	422	VAL	2.5
1	B	448	PHE	2.4
1	B	533	ASP	2.4
1	B	410	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	708	HIS	2.3
1	B	413	LYS	2.3
1	B	379	ASP	2.2
1	B	452	GLU	2.2
1	A	656	LYS	2.1
1	B	414	ASP	2.0
1	B	412	LYS	2.0
1	A	540	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SEP	B	695	10/11	0.86	0.17	89,94,100,101	0
1	SEP	A	676	10/11	0.95	0.09	57,60,64,65	0
1	SEP	A	695	10/11	0.96	0.10	40,45,53,54	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

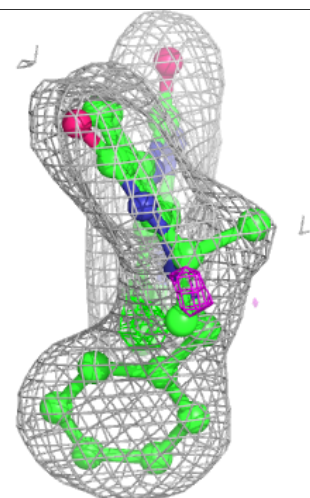
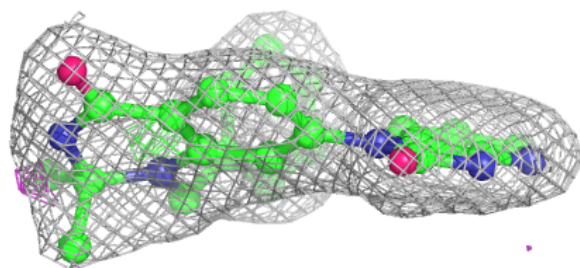
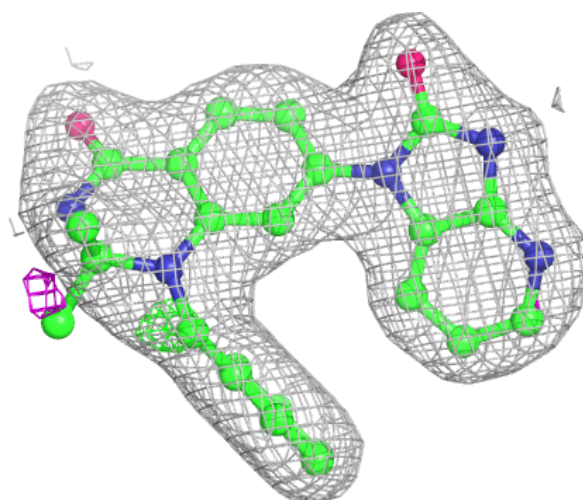
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	5VS	A	1001	30/30	0.97	0.14	35,38,43,47	0
2	5VS	B	1001	30/30	0.97	0.11	46,49,52,53	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.

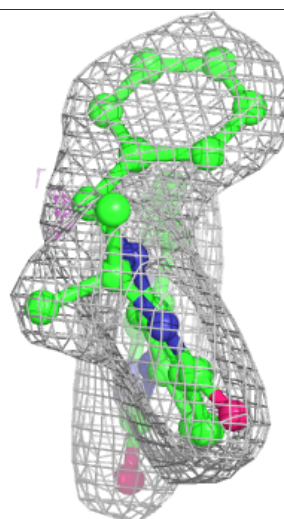
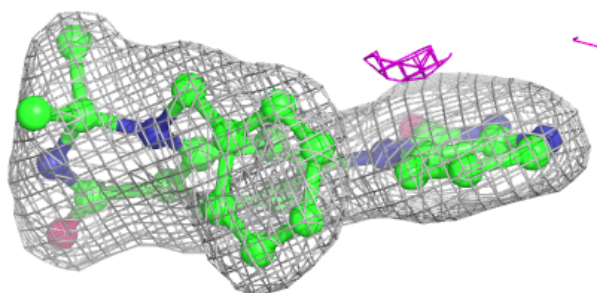
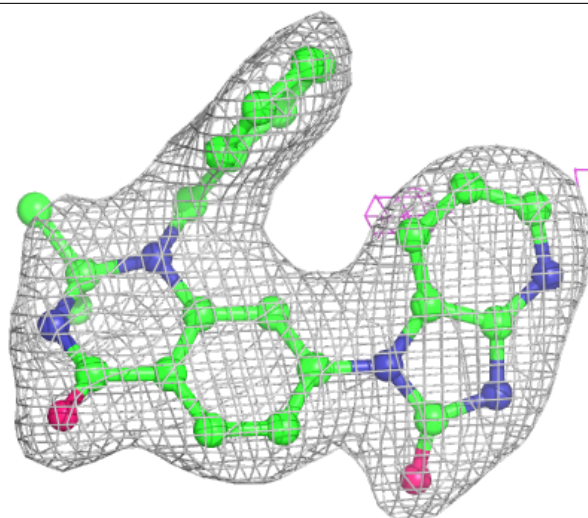
Electron density around 5VS A 1001:

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



Electron density around 5VS B 1001:

$2mF_o-DF_c$ (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.