



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:08 PM UTC

PDB ID : 6TFU / pdb_00006tfu
Title : Crystal Structure of EGFR T790M/V948R in Complex with Covalent Pyrrolopyrimidine 14d
Authors : Niggenaber, J.; Mueller, M.P.; Rauh, D.
Deposited on : 2019-11-14
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	:	4-5-2 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

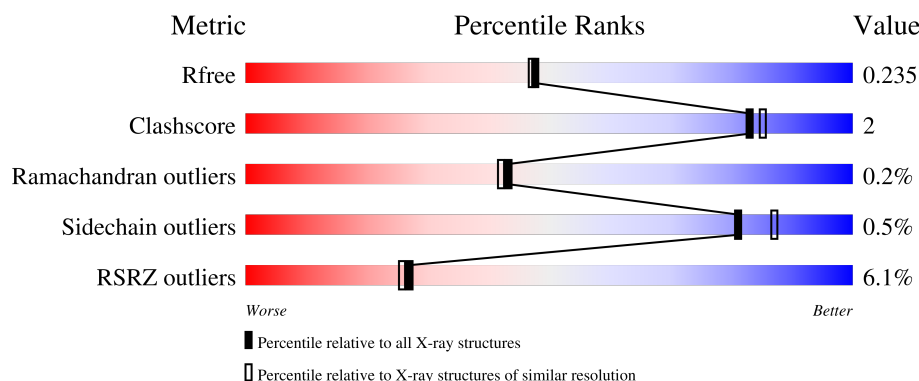
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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	333	4% 76% 5% 20%
1	B	333	6% 71% 7% 22%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	1102	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4200 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.

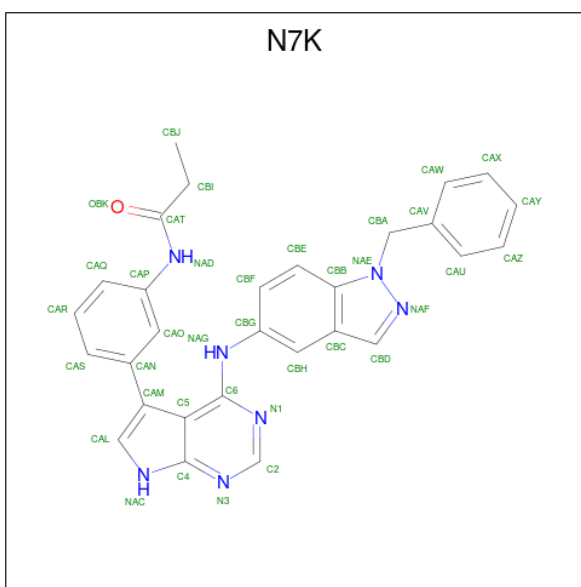
- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2041	1323	341	361	16			
1	B	260	Total	C	N	O	S	0	0	0
			1996	1294	342	345	15			

There are 14 discrepancies between the modelled and reference sequences:

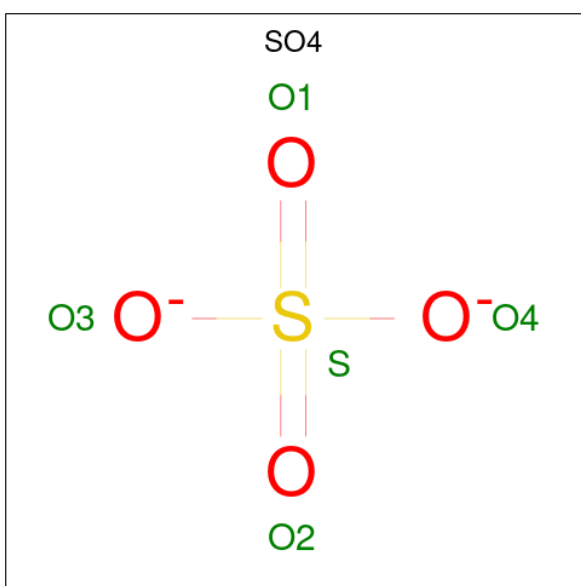
Chain	Residue	Modelled	Actual	Comment	Reference
A	690	GLY	-	expression tag	UNP P00533
A	691	SER	-	expression tag	UNP P00533
A	692	HIS	-	expression tag	UNP P00533
A	693	MET	-	expression tag	UNP P00533
A	694	ALA	-	expression tag	UNP P00533
A	790	MET	THR	engineered mutation	UNP P00533
A	948	ARG	VAL	engineered mutation	UNP P00533
B	690	GLY	-	expression tag	UNP P00533
B	691	SER	-	expression tag	UNP P00533
B	692	HIS	-	expression tag	UNP P00533
B	693	MET	-	expression tag	UNP P00533
B	694	ALA	-	expression tag	UNP P00533
B	790	MET	THR	engineered mutation	UNP P00533
B	948	ARG	VAL	engineered mutation	UNP P00533

- Molecule 2 is {N}-[3-[4-[[1-(phenylmethyl)indazol-5-yl]amino]-7 {H}-pyrrolo[2,3-d]pyrimidin-5-yl]phenyl]propanamide (CCD ID: N7K) (formula: C₂₉H₂₅N₇O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 37	C 29	N 7	O 1	0	0
2	B	1	Total 37	C 29	N 7	O 1	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	37	Total	O	0	0
			37	37		
4	B	32	Total	O	0	0
			32	32		

- Molecule 1: Epidermal growth factor receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.43Å 81.22Å 89.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.96 – 2.00 44.96 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.96-2.00) 100.0 (44.96-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.202 , 0.233 0.206 , 0.235	Depositor DCC
R_{free} test set	1926 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 59.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4200	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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: N7K, 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	0.15	0/2087	0.34	0/2845
1	B	0.15	0/2039	0.35	0/2769
All	All	0.15	0/4126	0.34	0/5614

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2041	0	1990	8	0
1	B	1996	0	1975	11	0
2	A	37	0	0	0	0
2	B	37	0	0	0	0
3	A	10	0	0	2	0
3	B	10	0	0	1	0
4	A	37	0	0	0	0
4	B	32	0	0	0	0
All	All	4200	0	3965	19	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 2.

All (19) 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:703:LEU:HD23	1:A:768:SER:HA	1.84	0.58
1:B:803:ARG:NH2	3:B:1103:SO4:O3	2.38	0.56
1:B:731:TRP:CE2	1:B:733:PRO:HG3	2.43	0.53
1:B:720:SER:HA	1:B:725:THR:HA	1.90	0.52
1:B:924:SER:O	1:B:928:GLU:HG2	2.12	0.50
1:B:835:HIS:CD2	1:B:856:PHE:HB3	2.47	0.49
1:B:943:VAL:HG22	1:B:971:MET:HE1	1.94	0.49
1:A:835:HIS:O	1:A:836:ARG:HB2	2.13	0.49
1:A:766:MET:HE3	1:A:777:LEU:HD22	1.96	0.47
3:A:1102:SO4:O3	1:B:846:LYS:NZ	2.48	0.46
1:B:714:LYS:HD3	1:B:727:TYR:CG	2.51	0.46
1:A:749:GLU:HB2	1:A:785:THR:HG22	1.98	0.45
1:A:846:LYS:NZ	3:A:1102:SO4:O1	2.46	0.45
1:A:730:LEU:HD23	1:A:730:LEU:HA	1.93	0.43
1:B:775:CYS:SG	1:B:790:MET:HE1	2.59	0.42
1:A:790:MET:HE1	1:A:844:LEU:HD12	2.01	0.42
1:A:926:ILE:HG23	1:A:931:GLU:HB2	2.01	0.41
1:B:882:ALA:HA	1:B:898:TRP:CD2	2.56	0.41
1:B:744:ILE:HG12	1:B:789:ILE:HD12	2.03	0.41

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	261/333 (78%)	256 (98%)	5 (2%)	0	100	100
1	B	252/333 (76%)	246 (98%)	5 (2%)	1 (0%)	30	27
All	All	513/666 (77%)	502 (98%)	10 (2%)	1 (0%)	43	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	748	ARG

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	210/291 (72%)	209 (100%)	1 (0%)	81	87
1	B	205/291 (70%)	204 (100%)	1 (0%)	81	87
All	All	415/582 (71%)	413 (100%)	2 (0%)	81	87

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	963	GLU
1	B	786	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	888	HIS

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

6 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
3	SO4	A	1102	-	4,4,4	0.26	0	6,6,6	0.11	0
3	SO4	B	1102	-	4,4,4	0.26	0	6,6,6	0.11	0
3	SO4	A	1103	-	4,4,4	0.27	0	6,6,6	0.16	0
3	SO4	B	1103	-	4,4,4	0.29	0	6,6,6	0.21	0
2	N7K	A	1101	1	42,42,42	3.88	17 (40%)	54,59,59	2.48	17 (31%)
2	N7K	B	1101	1	42,42,42	4.51	18 (42%)	54,59,59	2.58	16 (29%)

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	N7K	A	1101	1	-	4/18/18/18	0/6/6/6
2	N7K	B	1101	1	-	2/18/18/18	0/6/6/6

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	N7K	NAE-NAF	-21.63	1.22	1.37
2	A	1101	N7K	NAE-NAF	-16.43	1.26	1.37
2	B	1101	N7K	CAN-CAM	-10.12	1.34	1.49
2	A	1101	N7K	CAN-CAM	-9.95	1.34	1.49
2	B	1101	N7K	CBC-CBD	-8.04	1.33	1.42
2	A	1101	N7K	CBC-CBD	-7.69	1.33	1.42
2	A	1101	N7K	CBA-CAV	-5.69	1.41	1.51
2	B	1101	N7K	CBA-CAV	-5.54	1.41	1.51
2	B	1101	N7K	CBC-CBB	-5.53	1.33	1.41
2	A	1101	N7K	CBC-CBB	-5.12	1.33	1.41
2	B	1101	N7K	CAL-CAM	-4.44	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	N7K	C5-CAM	-4.33	1.35	1.46
2	A	1101	N7K	C5-CAM	-4.21	1.36	1.46
2	B	1101	N7K	CBH-CBC	-4.11	1.33	1.40
2	A	1101	N7K	CAL-CAM	-4.07	1.33	1.37
2	A	1101	N7K	CBH-CBC	-4.03	1.34	1.40
2	B	1101	N7K	C5-C4	-4.02	1.35	1.41
2	A	1101	N7K	C5-C4	-3.83	1.35	1.41
2	A	1101	N7K	CAP-NAD	-3.71	1.34	1.41
2	B	1101	N7K	CBE-CBB	-3.60	1.34	1.39
2	B	1101	N7K	CAP-NAD	-3.50	1.34	1.41
2	A	1101	N7K	CBE-CBB	-3.43	1.34	1.39
2	B	1101	N7K	C5-C6	-3.19	1.35	1.42
2	A	1101	N7K	C5-C6	-2.97	1.36	1.42
2	A	1101	N7K	C2-N1	2.91	1.39	1.33
2	B	1101	N7K	CAL-NAC	-2.91	1.32	1.37
2	B	1101	N7K	CBG-NAG	-2.62	1.35	1.40
2	B	1101	N7K	C2-N1	2.60	1.38	1.33
2	A	1101	N7K	C2-N3	2.56	1.38	1.33
2	B	1101	N7K	C2-N3	2.45	1.38	1.33
2	A	1101	N7K	C4-NAC	-2.41	1.33	1.36
2	A	1101	N7K	CAL-NAC	-2.35	1.33	1.37
2	B	1101	N7K	CBA-NAE	2.32	1.49	1.46
2	B	1101	N7K	C4-NAC	-2.19	1.33	1.36
2	A	1101	N7K	CBA-NAE	2.06	1.49	1.46

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1101	N7K	N3-C2-N1	-7.71	116.91	128.58
2	A	1101	N7K	N3-C2-N1	-7.40	117.39	128.58
2	B	1101	N7K	CBD-NAF-NAE	6.99	109.55	106.06
2	B	1101	N7K	C5-CAM-CAL	6.51	109.91	105.68
2	A	1101	N7K	C5-CAM-CAL	6.28	109.76	105.68
2	B	1101	N7K	CBC-CBD-NAF	-5.95	107.56	112.05
2	A	1101	N7K	CAP-NAD-CAT	-5.86	117.15	127.52
2	A	1101	N7K	CBC-CBD-NAF	-5.23	108.10	112.05
2	A	1101	N7K	CBD-NAF-NAE	5.20	108.65	106.06
2	B	1101	N7K	CBG-NAG-C6	-4.50	118.37	129.43
2	B	1101	N7K	CAP-NAD-CAT	-4.50	119.56	127.52
2	A	1101	N7K	CBG-NAG-C6	-4.44	118.53	129.43
2	B	1101	N7K	CAM-CAL-NAC	-4.04	107.85	110.46
2	A	1101	N7K	CAM-CAL-NAC	-4.02	107.87	110.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	N7K	C5-C4-N3	-3.95	122.43	126.64
2	B	1101	N7K	C5-C4-N3	-3.94	122.44	126.64
2	B	1101	N7K	CAN-CAM-CAL	-3.85	116.97	124.90
2	A	1101	N7K	CBB-CBC-CBD	3.60	106.52	104.48
2	A	1101	N7K	CBB-NAE-NAF	-3.46	109.70	111.26
2	B	1101	N7K	CBB-NAE-NAF	-3.16	109.83	111.26
2	B	1101	N7K	CAO-CAN-CAM	-3.14	116.66	120.74
2	A	1101	N7K	CAN-CAM-CAL	-3.12	118.47	124.90
2	A	1101	N7K	CBA-NAE-NAF	2.90	124.48	120.80
2	B	1101	N7K	CBA-NAE-NAF	2.74	124.27	120.80
2	A	1101	N7K	CBH-CBC-CBD	-2.67	131.55	135.46
2	A	1101	N7K	CBF-CBG-CBH	-2.65	116.45	119.66
2	B	1101	N7K	CBB-CBC-CBD	2.59	105.95	104.48
2	B	1101	N7K	CAS-CAN-CAM	2.52	124.14	120.98
2	A	1101	N7K	CBH-CBC-CBB	2.34	121.92	119.56
2	A	1101	N7K	CBC-CBH-CBG	2.29	121.50	119.00
2	A	1101	N7K	CBA-NAE-CBB	-2.22	125.82	128.71
2	B	1101	N7K	CBA-NAE-CBB	-2.17	125.89	128.71
2	B	1101	N7K	CBH-CBC-CBD	-2.04	132.48	135.46

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	N7K	CAQ-CAP-NAD-CAT
2	A	1101	N7K	CAO-CAP-NAD-CAT
2	B	1101	N7K	CAO-CAP-NAD-CAT
2	B	1101	N7K	CAQ-CAP-NAD-CAT
2	A	1101	N7K	CBF-CBG-NAG-C6
2	A	1101	N7K	CBH-CBG-NAG-C6

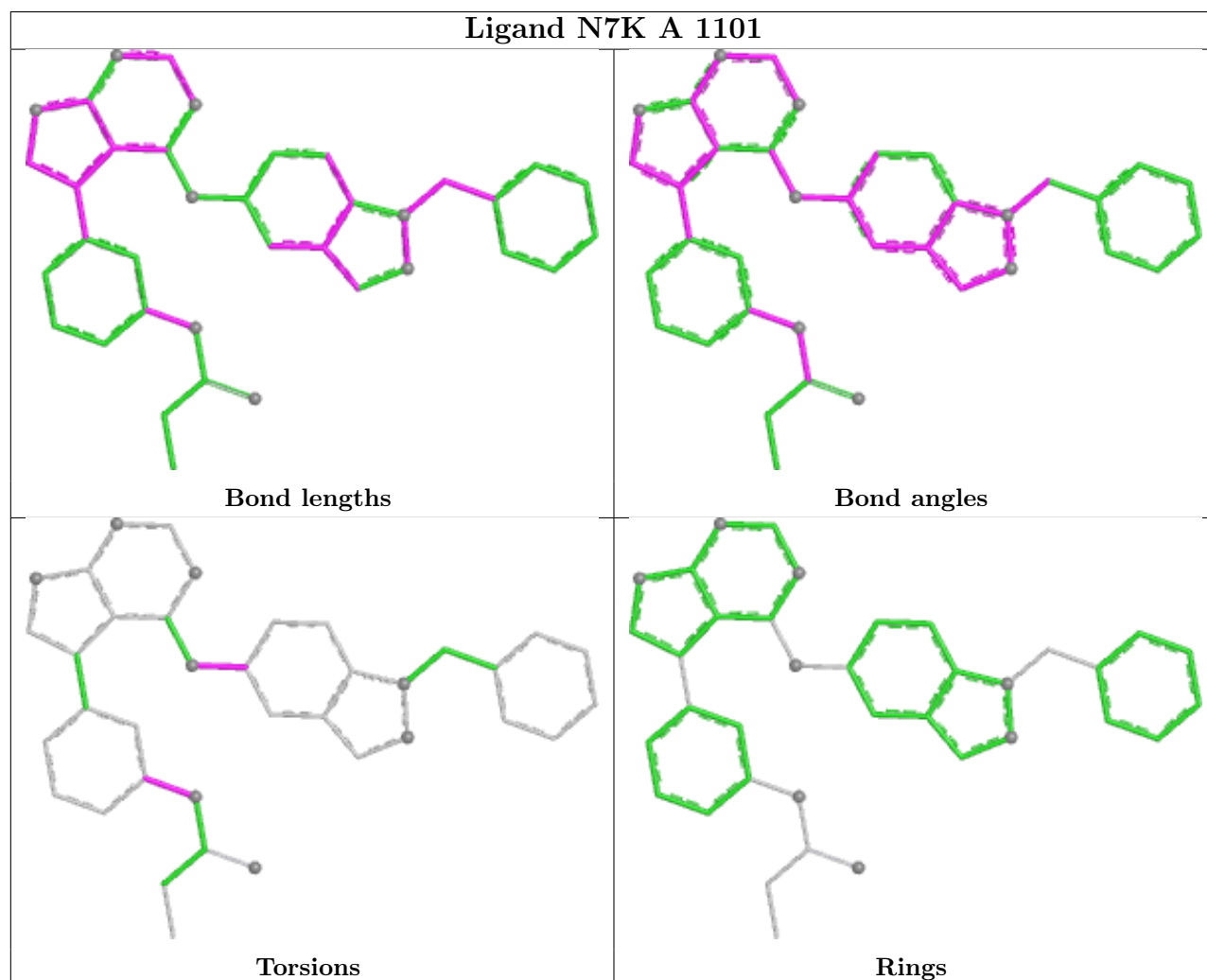
There are no ring outliers.

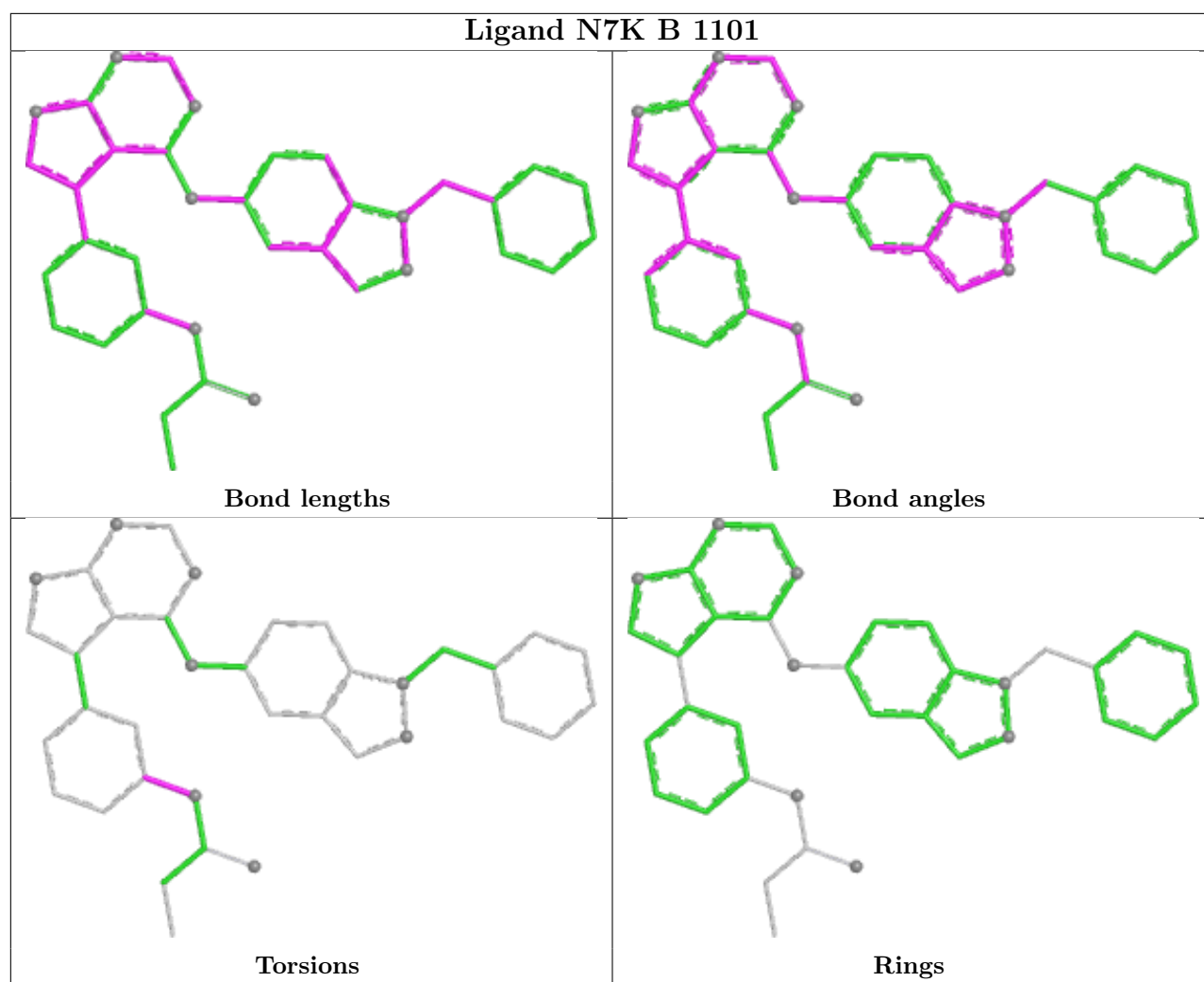
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1102	SO4	2	0
3	B	1103	SO4	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 [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	267/333 (80%)	0.46	12 (4%) 38 37	34, 51, 81, 102	0
1	B	260/333 (78%)	0.64	20 (7%) 19 18	34, 55, 91, 109	0
All	All	527/666 (79%)	0.55	32 (6%) 27 26	34, 52, 88, 109	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	723	PHE	5.0
1	B	859	ALA	4.7
1	B	759	ILE	4.2
1	A	700	ASN	4.2
1	A	751	THR	3.9
1	B	858	LEU	3.6
1	A	858	LEU	3.3
1	B	857	GLY	3.2
1	A	738	VAL	3.0
1	A	856	PHE	3.0
1	B	715	ILE	3.0
1	A	750	ALA	2.9
1	B	702	ALA	2.7
1	A	861	LEU	2.6
1	B	758	GLU	2.6
1	B	984	ASP	2.6
1	B	747	LEU	2.6
1	B	760	LEU	2.5
1	B	876	VAL	2.5
1	B	782	LEU	2.5
1	A	723	PHE	2.5
1	B	721	GLY	2.4
1	A	736	GLU	2.4
1	B	764	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	705	ARG	2.3
1	B	834	VAL	2.3
1	A	701	GLN	2.3
1	B	918	ILE	2.2
1	A	722	ALA	2.1
1	B	830	ASP	2.1
1	A	703	LEU	2.1
1	B	730	LEU	2.1

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](#)

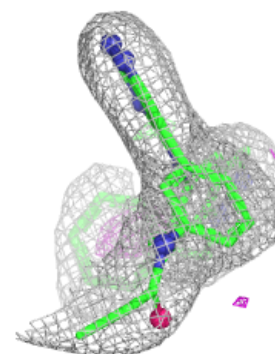
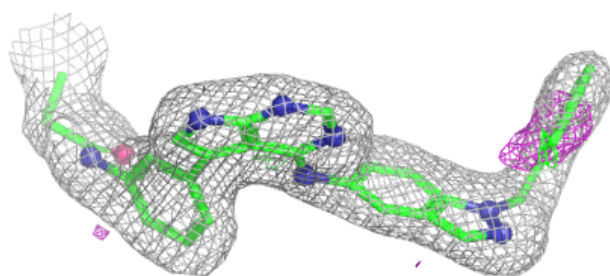
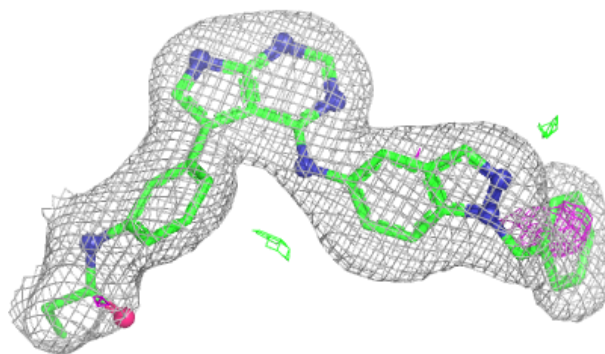
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	SO4	B	1103	5/5	0.81	0.09	109,110,112,112	0
3	SO4	A	1103	5/5	0.92	0.10	69,76,82,84	0
2	N7K	B	1101	37/37	0.92	0.10	40,50,63,76	0
2	N7K	A	1101	37/37	0.94	0.08	38,46,51,54	0
3	SO4	A	1102	5/5	0.94	0.09	70,71,75,79	0
3	SO4	B	1102	5/5	0.95	0.08	65,70,73,74	0

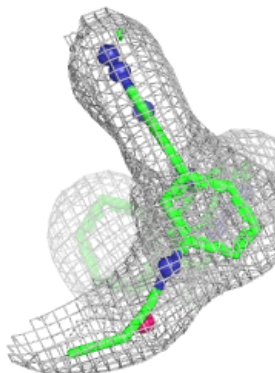
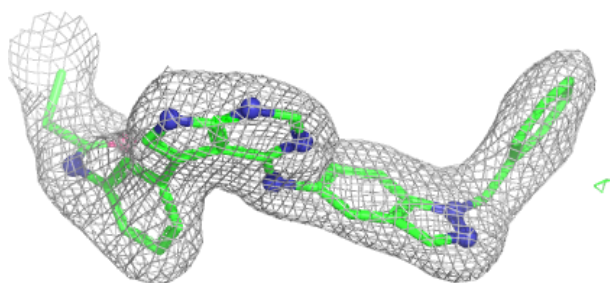
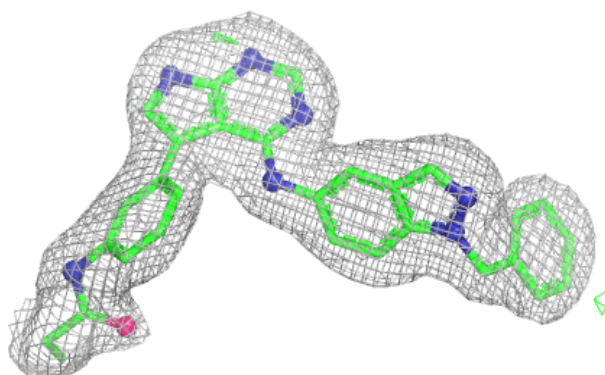
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around N7K B 1101:

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

**Electron density around N7K A 1101:**

$2mF_o-DF_c$ (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.