



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 03:05 PM UTC

PDB ID : 5IJ8 / pdb_00005ij8
Title : Structure of the primary oncogenic mutant Y641N Hs/AcPRC2 in complex with a pyridone inhibitor
Authors : Gajiwala, K.S.; Brooun, A.; Deng, Y.-L.; Liu, W.
Deposited on : 2016-03-01
Resolution : 2.99 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

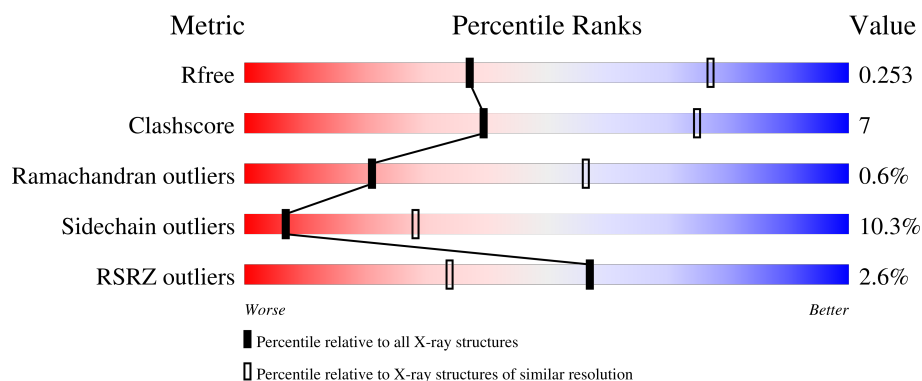
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	643	
1	B	643	
2	E	362	
2	F	362	
3	S	191	

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Mol	Chain	Length	Quality of chain
3	T	191	3% 45% 19% • 33%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15580 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 Enhancer of Zeste Homolog 2 (EZH2), Histone-lysine N-methyltransferase EZH2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3771	2363	680	689	39			
1	B	466	Total	C	N	O	S	0	0	0
			3779	2366	681	693	39			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	PRO	deletion	UNP Q15910
A	?	-	PRO	deletion	UNP Q15910
A	?	-	ARG	deletion	UNP Q15910
A	?	-	LYS	deletion	UNP Q15910
A	492	LEU	LYS	linker	UNP Q15910
A	493	GLY	LYS	linker	UNP Q15910
A	494	GLY	ARG	linker	UNP Q15910
A	495	GLY	LYS	linker	UNP Q15910
A	496	GLY	HIS	linker	UNP Q15910
A	497	SER	ARG	linker	UNP Q15910
A	498	GLY	LEU	linker	UNP Q15910
A	499	GLY	TRP	linker	UNP Q15910
A	500	GLY	ALA	linker	UNP Q15910
A	501	GLY	ALA	linker	UNP Q15910
A	502	SER	HIS	linker	UNP Q15910
A	503	GLY	CYS	linker	UNP Q15910
A	504	GLY	ARG	linker	UNP Q15910
A	505	GLY	LYS	linker	UNP Q15910
A	506	GLY	ILE	linker	UNP Q15910
A	507	SER	GLN	linker	UNP Q15910
A	508	ALA	LEU	linker	UNP Q15910
A	509	ALA	LYS	linker	UNP Q15910
A	510	ALA	LYS	linker	UNP Q15910
A	532	ASN	SER	linker	UNP Q15910

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Chain	Residue	Modelled	Actual	Comment	Reference
A	641	ASN	TYR	engineered mutation	UNP Q15910
B	?	-	PRO	deletion	UNP Q15910
B	?	-	PRO	deletion	UNP Q15910
B	?	-	ARG	deletion	UNP Q15910
B	?	-	LYS	deletion	UNP Q15910
B	492	LEU	LYS	linker	UNP Q15910
B	493	GLY	LYS	linker	UNP Q15910
B	494	GLY	ARG	linker	UNP Q15910
B	495	GLY	LYS	linker	UNP Q15910
B	496	GLY	HIS	linker	UNP Q15910
B	497	SER	ARG	linker	UNP Q15910
B	498	GLY	LEU	linker	UNP Q15910
B	499	GLY	TRP	linker	UNP Q15910
B	500	GLY	ALA	linker	UNP Q15910
B	501	GLY	ALA	linker	UNP Q15910
B	502	SER	HIS	linker	UNP Q15910
B	503	GLY	CYS	linker	UNP Q15910
B	504	GLY	ARG	linker	UNP Q15910
B	505	GLY	LYS	linker	UNP Q15910
B	506	GLY	ILE	linker	UNP Q15910
B	507	SER	GLN	linker	UNP Q15910
B	508	ALA	LEU	linker	UNP Q15910
B	509	ALA	LYS	linker	UNP Q15910
B	510	ALA	LYS	linker	UNP Q15910
B	532	ASN	SER	linker	UNP Q15910
B	641	ASN	TYR	engineered mutation	UNP Q15910

- Molecule 2 is a protein called Polycomb protein EED.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	358	Total	C	N	O	S	0	0	0
			2898	1834	509	534	21			
2	F	359	Total	C	N	O	S	0	0	0
			2906	1840	510	535	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	66	MET	-	initiating methionine	UNP O75530
F	66	MET	-	initiating methionine	UNP O75530

- Molecule 3 is a protein called Polycomb protein SUZ12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	S	128	Total	C	N	O	S	0	0	0
			1070	675	185	198	12			
3	T	128	Total	C	N	O	S	0	0	0
			1070	675	185	198	12			

There are 22 discrepancies between the modelled and reference sequences:

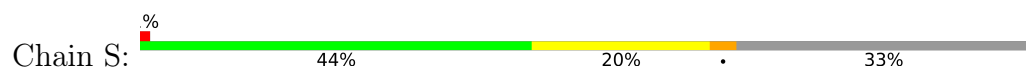
Chain	Residue	Modelled	Actual	Comment	Reference
S	535	MET	-	initiating methionine	UNP Q15022
S	536	ASP	-	expression tag	UNP Q15022
S	537	TYR	-	expression tag	UNP Q15022
S	538	LYS	-	expression tag	UNP Q15022
S	539	ASP	-	expression tag	UNP Q15022
S	540	ASP	-	expression tag	UNP Q15022
S	541	ASP	-	expression tag	UNP Q15022
S	542	ASP	-	expression tag	UNP Q15022
S	543	LYS	-	expression tag	UNP Q15022
S	544	GLY	-	expression tag	UNP Q15022
S	583	ASP	SER	engineered mutation	UNP Q15022
T	535	MET	-	initiating methionine	UNP Q15022
T	536	ASP	-	expression tag	UNP Q15022
T	537	TYR	-	expression tag	UNP Q15022
T	538	LYS	-	expression tag	UNP Q15022
T	539	ASP	-	expression tag	UNP Q15022
T	540	ASP	-	expression tag	UNP Q15022
T	541	ASP	-	expression tag	UNP Q15022
T	542	ASP	-	expression tag	UNP Q15022
T	543	LYS	-	expression tag	UNP Q15022
T	544	GLY	-	expression tag	UNP Q15022
T	583	ASP	SER	engineered mutation	UNP Q15022

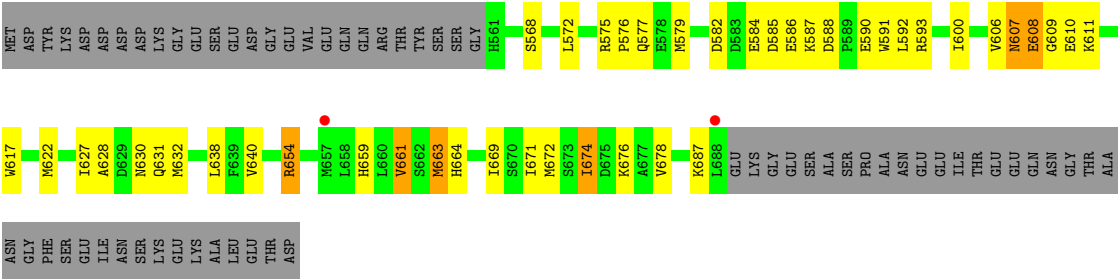
- Molecule 4 is 5,8-dichloro-2-[(4-ethyl-6-methyl-2-oxo-1,2-dihydropyridin-3-yl)methyl]-7-(1-[(2R)-2-hydroxypropanoyl]piperidin-4-yl}oxy)-3,4-dihydroisoquinolin-1(2H)-one (CCD ID: 6BN) (formula: C₂₆H₃₁Cl₂N₃O₅).



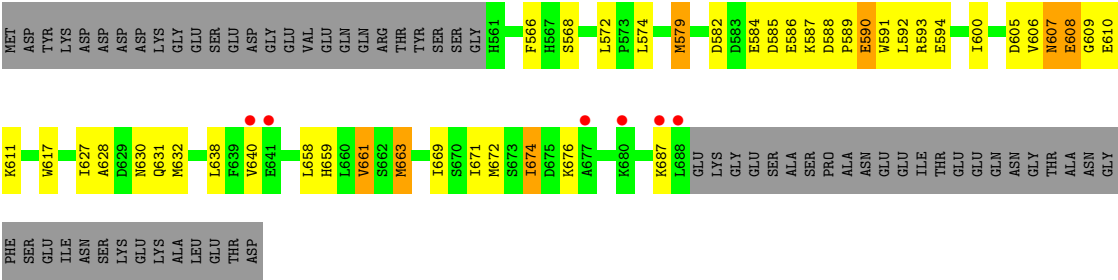
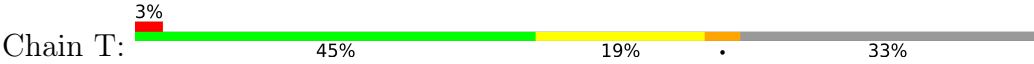
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	7	Total Zn 7 7	0	0
5	B	7	Total Zn 7 7	0	0





● Molecule 3: Polycomb protein SUZ12



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.33Å 115.13Å 153.94Å 90.00° 103.35° 90.00°	Depositor
Resolution (Å)	91.28 – 2.99 91.28 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.7 (91.28-2.99) 99.7 (91.28-2.99)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.188 , 0.240 0.198 , 0.253	Depositor DCC
R_{free} test set	2497 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	64.4	Xtriage
Anisotropy	0.532	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 73.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15580	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.32% 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: 6BN, ZN

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	0.88	0/3851	1.39	38/5177 (0.7%)
1	B	0.87	0/3857	1.40	25/5185 (0.5%)
2	E	0.85	1/2972 (0.0%)	1.25	10/4026 (0.2%)
2	F	0.85	0/2980	1.28	9/4037 (0.2%)
3	S	0.86	0/1091	1.48	8/1464 (0.5%)
3	T	0.86	0/1091	1.50	6/1464 (0.4%)
All	All	0.86	1/15842 (0.0%)	1.36	96/21353 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	242	MET	SD-CE	5.68	1.93	1.79

The worst 5 of 96 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	675	VAL	N-CA-CB	9.87	121.48	110.72
1	A	675	VAL	N-CA-CB	9.44	121.00	110.72
2	F	356	MET	CA-C-N	9.04	134.40	120.75
2	F	356	MET	C-N-CA	9.04	134.40	120.75
1	B	447	TYR	N-CA-C	-6.87	100.36	110.52

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	3771	0	3661	41	0
1	B	3779	0	3674	48	0
2	E	2898	0	2812	49	0
2	F	2906	0	2823	41	0
3	S	1070	0	1052	26	0
3	T	1070	0	1052	27	0
4	A	36	0	0	0	0
4	B	36	0	0	0	0
5	A	7	0	0	0	0
5	B	7	0	0	0	0
All	All	15580	0	15074	202	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 7.

The worst 5 of 202 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:B:16:ARG:HD2	1:B:230:MET:HE3	1.58	0.85
3:S:628:ALA:H	3:S:631:GLN:HE21	1.26	0.83
3:T:628:ALA:H	3:T:631:GLN:HE21	1.26	0.82
1:B:670:ASN:HD22	1:B:672:ASP:H	1.28	0.79
3:T:627:ILE:H	3:T:631:GLN:NE2	1.81	0.79

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	450/643 (70%)	413 (92%)	33 (7%)	4 (1%)	14 48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	452/643 (70%)	418 (92%)	32 (7%)	2 (0%)	30	65
2	E	356/362 (98%)	343 (96%)	11 (3%)	2 (1%)	21	56
2	F	357/362 (99%)	344 (96%)	11 (3%)	2 (1%)	21	56
3	S	126/191 (66%)	112 (89%)	13 (10%)	1 (1%)	16	50
3	T	126/191 (66%)	116 (92%)	9 (7%)	1 (1%)	16	50
All	All	1867/2392 (78%)	1746 (94%)	109 (6%)	12 (1%)	21	56

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	593	HIS
1	B	593	HIS
2	F	380	GLU
1	A	245	LYS
1	B	460	THR

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	419/560 (75%)	372 (89%)	47 (11%)	6	24
1	B	421/560 (75%)	378 (90%)	43 (10%)	7	29
2	E	321/325 (99%)	294 (92%)	27 (8%)	10	37
2	F	322/325 (99%)	294 (91%)	28 (9%)	9	35
3	S	122/175 (70%)	105 (86%)	17 (14%)	3	17
3	T	122/175 (70%)	106 (87%)	16 (13%)	4	19
All	All	1727/2120 (82%)	1549 (90%)	178 (10%)	7	28

5 of 178 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	408	SER

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Mol	Chain	Res	Type
2	F	408	SER
2	F	166	MET
2	F	328	LYS
3	S	587	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 59 such sidechains are listed below:

Mol	Chain	Res	Type
2	E	175	HIS
3	T	642	ASN
2	F	86	GLN
3	T	631	GLN
3	S	642	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](#)

Of 16 ligands modelled in this entry, 14 are monoatomic - leaving 2 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
4	6BN	A	9001	-	38,39,39	0.43	0	46,57,57	0.91	2 (4%)
4	6BN	B	9001	-	38,39,39	0.34	0	46,57,57	0.87	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
4	6BN	A	9001	-	-	2/18/41/41	0/4/4/4
4	6BN	B	9001	-	-	2/18/41/41	0/4/4/4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	9001	6BN	C15-C14-C13	-2.69	117.28	120.26
4	A	9001	6BN	C4-C5-C8	-2.41	116.31	119.58
4	B	9001	6BN	C15-C14-C13	-2.39	117.61	120.26
4	B	9001	6BN	C4-C5-C8	-2.38	116.35	119.58

There are no chirality outliers.

All (4) torsion outliers are listed below:

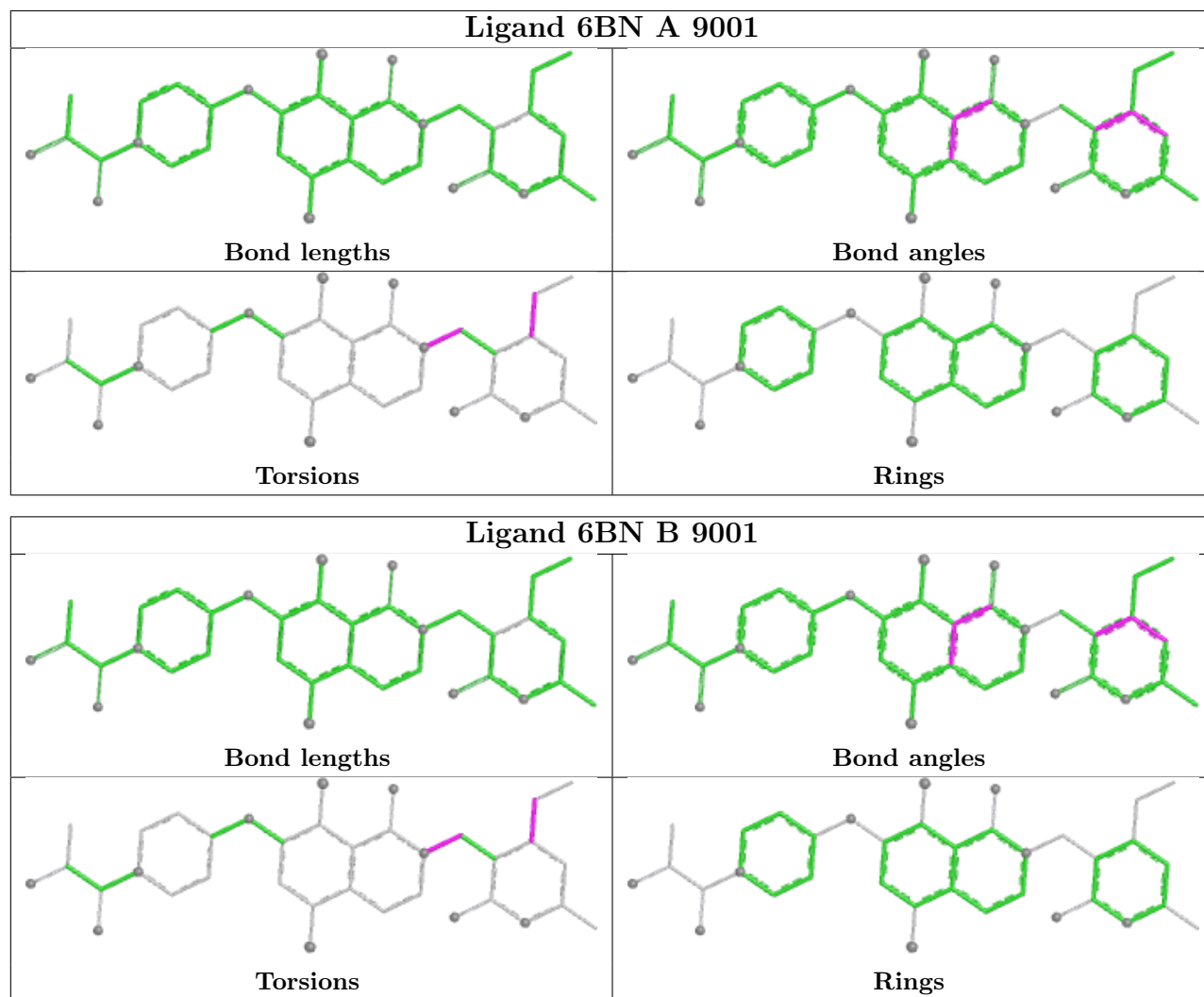
Mol	Chain	Res	Type	Atoms
4	A	9001	6BN	C13-C14-C21-C22
4	B	9001	6BN	C13-C14-C21-C22
4	A	9001	6BN	C13-C12-N9-C8
4	B	9001	6BN	C13-C12-N9-C8

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 [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	464/643 (72%)	0.19	17 (3%) 45 25	39, 66, 109, 139	0
1	B	466/643 (72%)	0.28	19 (4%) 41 23	41, 72, 113, 125	0
2	E	358/362 (98%)	-0.10	3 (0%) 82 64	37, 59, 87, 127	0
2	F	359/362 (99%)	-0.18	3 (0%) 82 64	35, 56, 87, 120	0
3	S	128/191 (67%)	0.45	2 (1%) 70 47	47, 83, 111, 117	0
3	T	128/191 (67%)	0.55	6 (4%) 36 19	41, 80, 112, 128	0
All	All	1903/2392 (79%)	0.13	50 (2%) 57 34	35, 65, 108, 139	0

The worst 5 of 50 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	219	PRO	4.3
1	B	301	ALA	3.8
1	B	219	PRO	3.6
3	T	688	LEU	3.6
1	A	733	ALA	3.5

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

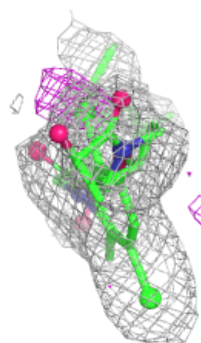
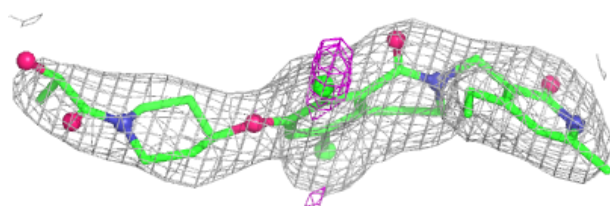
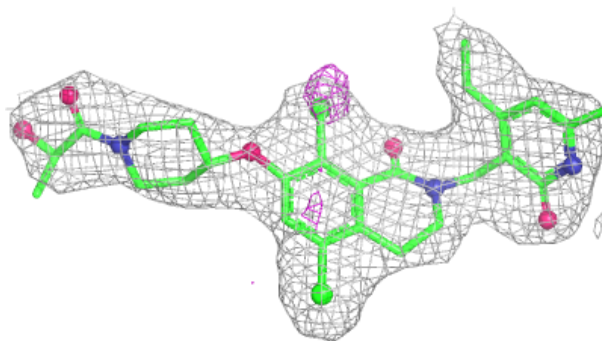
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	ZN	A	9005	1/1	0.93	0.21	300,300,300,300	1
4	6BN	B	9001	36/36	0.94	0.09	29,50,68,69	0
4	6BN	A	9001	36/36	0.95	0.08	18,46,55,56	0
5	ZN	A	9004	1/1	0.99	0.03	45,45,45,45	1
5	ZN	B	9002	1/1	0.99	0.07	53,53,53,53	1
5	ZN	B	9003	1/1	0.99	0.05	61,61,61,61	1
5	ZN	B	9004	1/1	0.99	0.08	59,59,59,59	1
5	ZN	B	9007	1/1	0.99	0.05	65,65,65,65	1
5	ZN	A	9008	1/1	1.00	0.04	35,35,35,35	1
5	ZN	A	9002	1/1	1.00	0.06	59,59,59,59	1
5	ZN	A	9003	1/1	1.00	0.03	45,45,45,45	1
5	ZN	A	9006	1/1	1.00	0.06	37,37,37,37	1
5	ZN	B	9005	1/1	1.00	0.07	42,42,42,42	1
5	ZN	B	9006	1/1	1.00	0.07	53,53,53,53	1
5	ZN	A	9007	1/1	1.00	0.06	39,39,39,39	1
5	ZN	B	9008	1/1	1.00	0.07	47,47,47,47	1

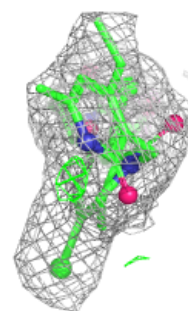
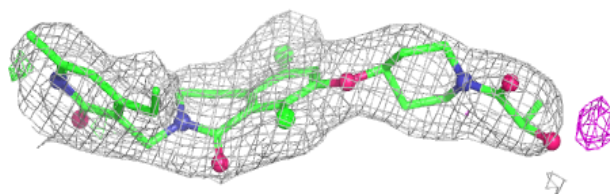
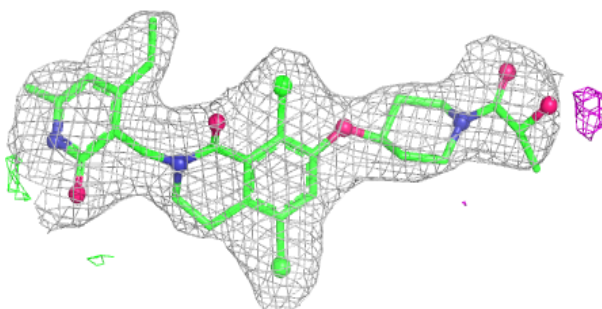
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6BN B 9001:

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

**Electron density around 6BN A 9001:**

$2mF_o-DF_c$ (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.