



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:41 PM UTC

PDB ID : 5J4Y / pdb_00005j4y
Title : The crystal structure of N-(4-(2-(thiazolo[5,4-c]pyridin-2-yl)phenoxy)phenyl)acetamide bound to JCV Helicase
Authors : Ter Haar, E.
Deposited on : 2016-04-01
Resolution : 2.59 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

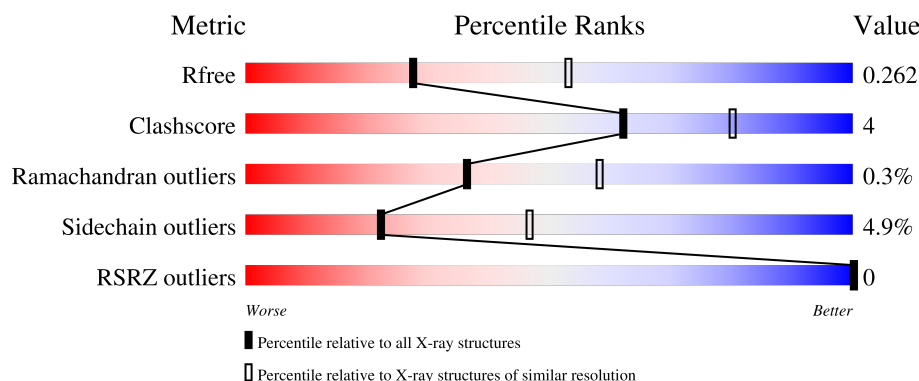
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	372	 81% 15% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3019 atoms, of which 12 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 Large T antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2868	1840	491	514	23			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	257	GLY	-	expression tag	UNP P03072
A	258	SER	-	expression tag	UNP P03072
A	259	HIS	-	expression tag	UNP P03072
A	260	MET	-	expression tag	UNP P03072
A	280	ASP	GLU	engineered mutation	UNP P03072
A	295	ASN	ASP	engineered mutation	UNP P03072
A	299	ALA	ASN	engineered mutation	UNP P03072
A	301	ALA	GLN	engineered mutation	UNP P03072
A	302	ALA	GLN	engineered mutation	UNP P03072
A	304	ALA	LYS	engineered mutation	UNP P03072
A	305	ALA	LYS	engineered mutation	UNP P03072
A	307	ALA	GLU	engineered mutation	UNP P03072
A	308	ALA	LYS	engineered mutation	UNP P03072
A	309	ALA	LYS	engineered mutation	UNP P03072
A	354	LEU	ILE	engineered mutation	UNP P03072
A	408	GLU	ASP	engineered mutation	UNP P03072
A	624	ALA	ARG	engineered mutation	UNP P03072

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	S	0	0
			24	6	12	1	4	1		

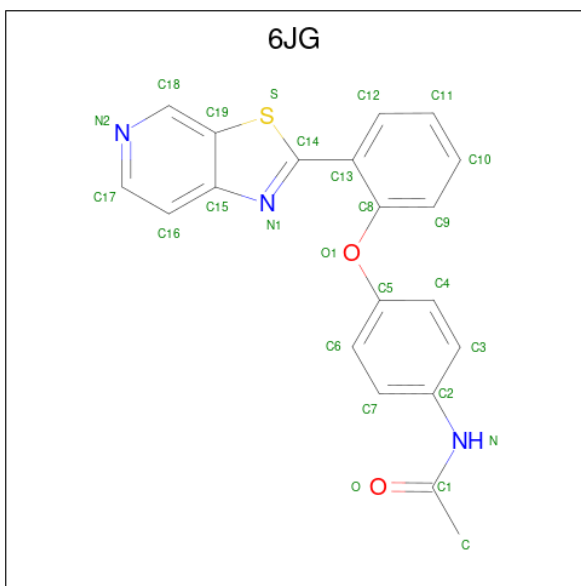
- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is N-{4-[2-([1,3]thiazolo[5,4-c]pyridin-2-yl)phenoxy]phenyl}acetamide (CCD ID:

6JG) (formula: C₂₀H₁₅N₃O₂S).



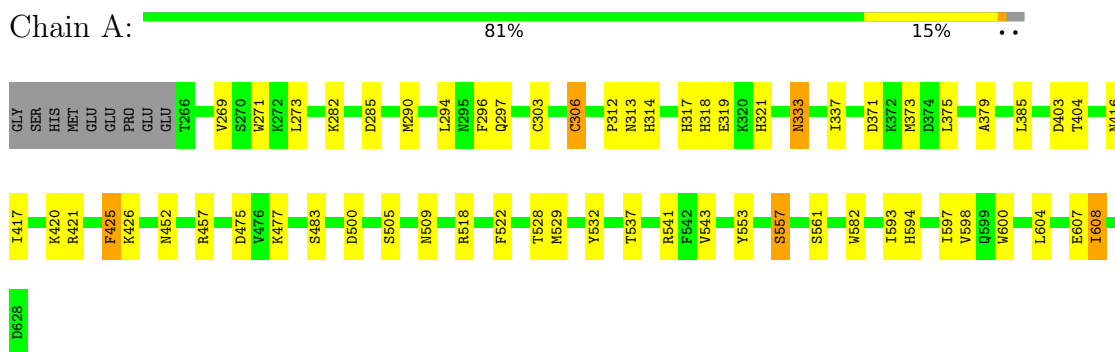
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			26	20	3	2	1		

- Molecule 6 is water.

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

i

- Molecule 1: Large T antigen



4 Data and refinement statistics

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	109.25Å 109.25Å 66.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	54.62 – 2.59 54.62 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.4 (54.62-2.59) 98.4 (54.62-2.59)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.58Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.170 , 0.238 0.182 , 0.262	Depositor DCC
R_{free} test set	699 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	61.5	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.175 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3019	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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: ZN, 6JG, MES, 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	0.87	0/2929	1.42	18/3962 (0.5%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	416	ASN	CA-CB-CG	6.48	119.08	112.60
1	A	425	PHE	CA-CB-CG	6.47	120.27	113.80
1	A	371	ASP	CA-C-N	5.86	128.13	120.28
1	A	371	ASP	C-N-CA	5.86	128.13	120.28
1	A	296	PHE	CA-CB-CG	5.81	119.61	113.80
1	A	505	SER	N-CA-C	-5.80	105.88	113.12
1	A	285	ASP	CA-CB-CG	5.70	118.30	112.60
1	A	557	SER	CA-C-N	5.55	127.66	120.44
1	A	557	SER	C-N-CA	5.55	127.66	120.44
1	A	271	TRP	CA-C-N	5.54	127.70	120.28
1	A	271	TRP	C-N-CA	5.54	127.70	120.28
1	A	561	SER	CA-C-N	5.37	127.73	120.38
1	A	561	SER	C-N-CA	5.37	127.73	120.38
1	A	403	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	333	ASN	CA-C-N	5.08	127.05	120.44
1	A	333	ASN	C-N-CA	5.08	127.05	120.44
1	A	457	ARG	CA-C-N	5.08	127.09	120.28
1	A	457	ARG	C-N-CA	5.08	127.09	120.28

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	2868	0	2866	23	0
2	A	1	0	0	0	0
3	A	12	12	13	0	0
4	A	10	0	0	0	0
5	A	26	0	0	1	0
6	A	90	0	0	0	0
All	All	3007	12	2879	24	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 4.

All (24) 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:452:ASN:HD21	1:A:477:LYS:H	1.39	0.70
1:A:317:HIS:HB3	1:A:321:HIS:HD2	1.67	0.59
1:A:385:LEU:HD23	1:A:607:GLU:HG3	1.89	0.53
1:A:303:CYS:HB3	1:A:306:CYS:HB2	1.91	0.52
1:A:417:ILE:O	1:A:421:ARG:HG3	2.11	0.50
1:A:314:HIS:O	1:A:318:HIS:HB2	2.12	0.50
1:A:426:LYS:HD2	1:A:529:MET:HE2	1.94	0.50
1:A:313:ASN:HA	1:A:317:HIS:CD2	2.49	0.48
1:A:425:PHE:HB2	1:A:528:THR:HG22	1.97	0.47
1:A:594:HIS:O	1:A:598:VAL:HG23	2.15	0.47
1:A:593:ILE:O	1:A:597:ILE:HG12	2.15	0.46
5:A:705:6JG:S	5:A:705:6JG:O1	2.73	0.46
1:A:553:TYR:O	1:A:557:SER:HB2	2.14	0.46
1:A:317:HIS:HB3	1:A:321:HIS:CD2	2.49	0.46
1:A:333:ASN:O	1:A:337:ILE:HG13	2.16	0.45
1:A:312:PRO:HD3	1:A:375:LEU:HD22	1.99	0.44
1:A:373:MET:HB3	1:A:600:TRP:CZ2	2.52	0.44
1:A:420:LYS:HA	1:A:543:VAL:HB	2.00	0.44
1:A:582:TRP:CH2	1:A:608:ILE:HG13	2.53	0.43
1:A:421:ARG:NH2	1:A:522:PHE:O	2.53	0.42
1:A:593:ILE:HD12	1:A:597:ILE:HD11	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:529:MET:HE3	1:A:532:TYR:HB2	2.01	0.41
1:A:294:LEU:O	1:A:297:GLN:HG2	2.20	0.41
1:A:385:LEU:HD21	1:A:604:LEU:HD23	2.03	0.41

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	361/372 (97%)	346 (96%)	14 (4%)	1 (0%)	36 58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	379	ALA

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	306/322 (95%)	291 (95%)	15 (5%)	22 47

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	269	VAL
1	A	273	LEU
1	A	282	LYS
1	A	290	MET
1	A	306	CYS
1	A	319	GLU
1	A	404	THR
1	A	475	ASP
1	A	483	SER
1	A	500	ASP
1	A	509	ASN
1	A	518	ARG
1	A	537	THR
1	A	541	ARG
1	A	608	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	297	GLN
1	A	317	HIS
1	A	326	GLN
1	A	333	ASN
1	A	382	ASN
1	A	452	ASN
1	A	509	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	SO4	A	703	-	4,4,4	0.21	0	6,6,6	0.24	0
3	MES	A	702	-	12,12,12	1.21	1 (8%)	15,16,16	0.59	0
5	6JG	A	705	-	29,29,29	0.24	0	36,40,40	0.49	1 (2%)
4	SO4	A	704	-	4,4,4	0.24	0	6,6,6	0.24	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MES	A	702	-	-	4/6/14/14	0/1/1/1
5	6JG	A	705	-	-	0/12/12/12	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	MES	C8-S	-4.06	1.71	1.77

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	705	6JG	C8-O1-C5	2.18	123.24	117.95

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	MES	C8-C7-N4-C3

Continued on next page...

Continued from previous page...

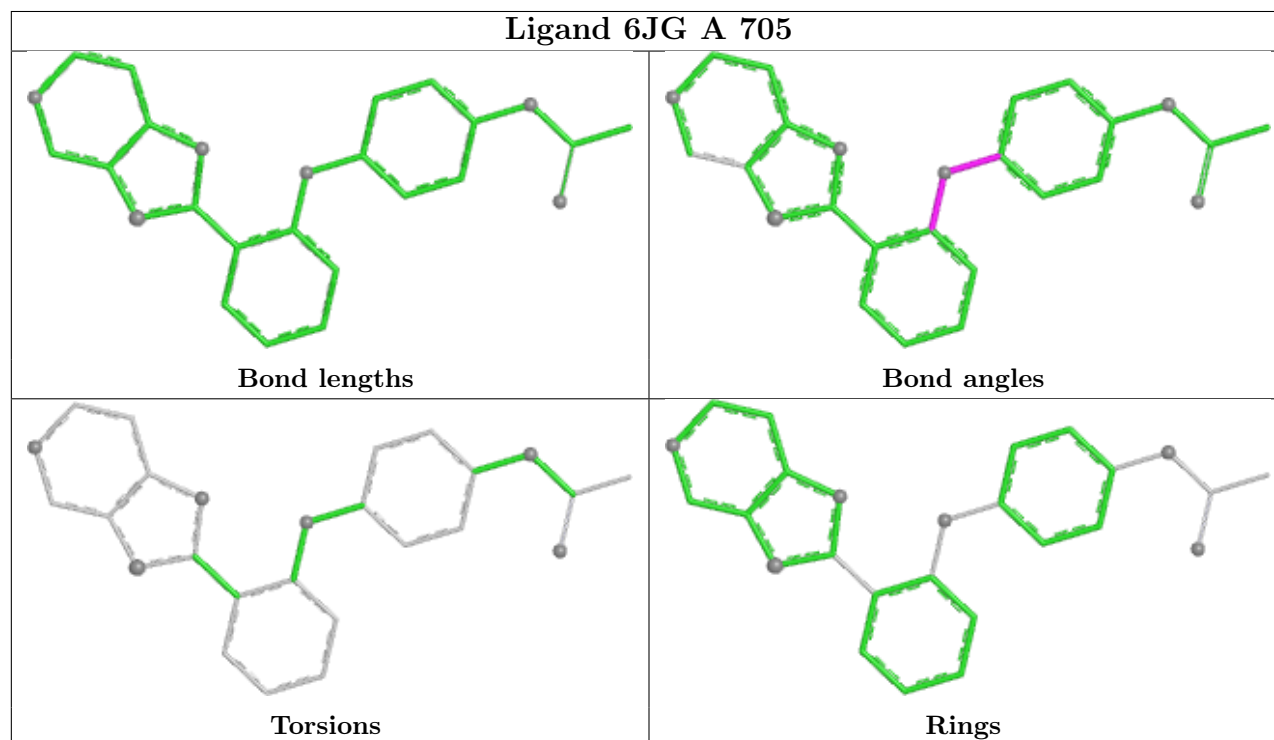
Mol	Chain	Res	Type	Atoms
3	A	702	MES	C7-C8-S-O3S
3	A	702	MES	C7-C8-S-O1S
3	A	702	MES	C7-C8-S-O2S

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	705	6JG	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 [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	363/372 (97%)	-1.25	0 100 100	41, 71, 118, 163	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

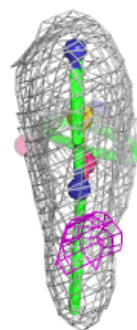
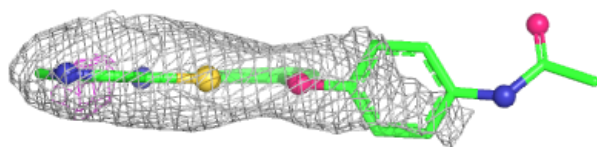
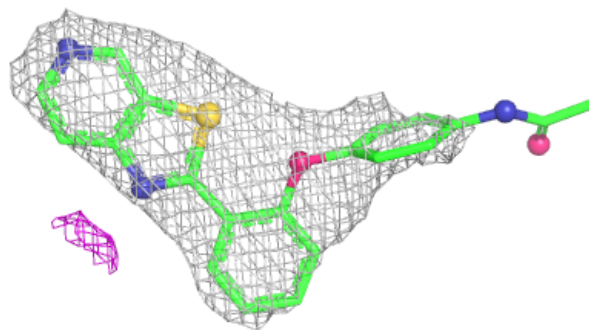
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MES	A	702	12/12	0.99	0.03	67,70,71,74	0
4	SO4	A	703	5/5	0.99	0.04	101,103,105,105	0
4	SO4	A	704	5/5	0.99	0.04	109,109,109,110	0
5	6JG	A	705	26/26	0.99	0.05	79,89,109,111	0
2	ZN	A	701	1/1	1.00	0.02	142,142,142,142	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 6JG A 705:

$2mF_o - DF_c$ (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.