



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 06:14 AM UTC

PDB ID : 5PZL / pdb_00005pzl
Title : CRYSTAL STRUCTURE OF THE HEPATITIS C VIRUS NS5B RNA-DEPENDENT RNA POLYMERASE IN COMPLEX WITH 2-({3-[1-(2-CYCLOPROPYLETHYL)-6-FLUORO-4-HYDROXY-2-OXO-1,2-DIHYDROQUINOLIN-3-YL]-1,1-DIOXO-1,4-DIHYDRO-1LAMBDA 6 ,2,4-BENZOTHIADIAZIN-7-YL}OXY)ACETAMIDE
Authors : Sheriff, S.
Deposited on : 2017-02-27
Resolution : 2.06 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)

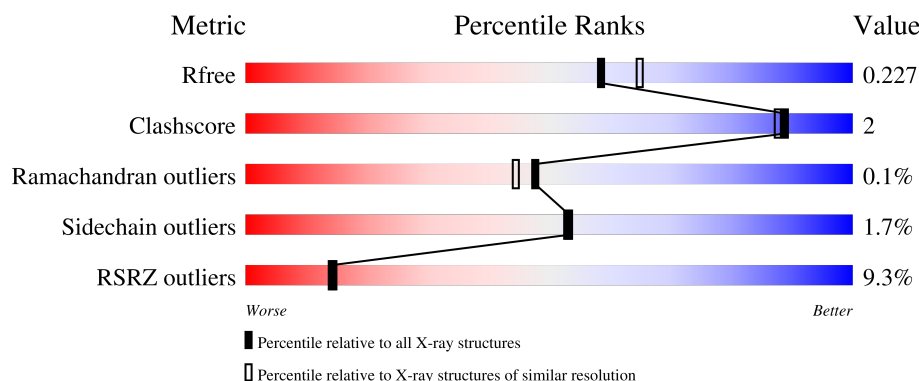
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	574	9% 90% 6% • •
1	B	574	9% 89% 5% • 5%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9492 atoms, of which 118 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.

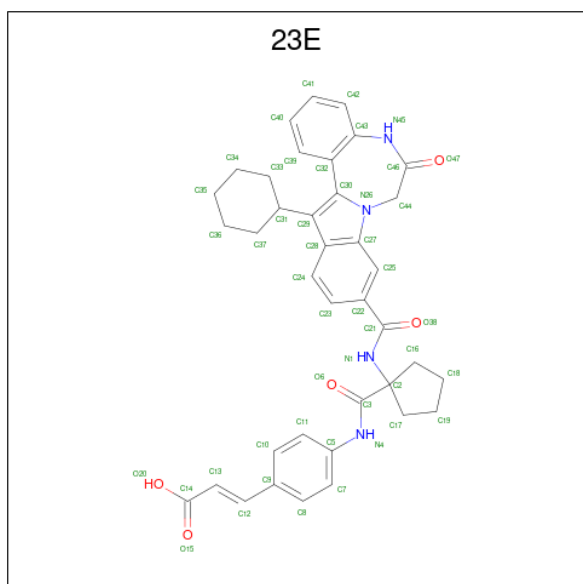
- Molecule 1 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	0
			4240	2675	746	786	33			
1	B	545	Total	C	N	O	S	0	0	0
			4182	2635	738	777	32			

There are 2 discrepancies between the modelled and reference sequences:

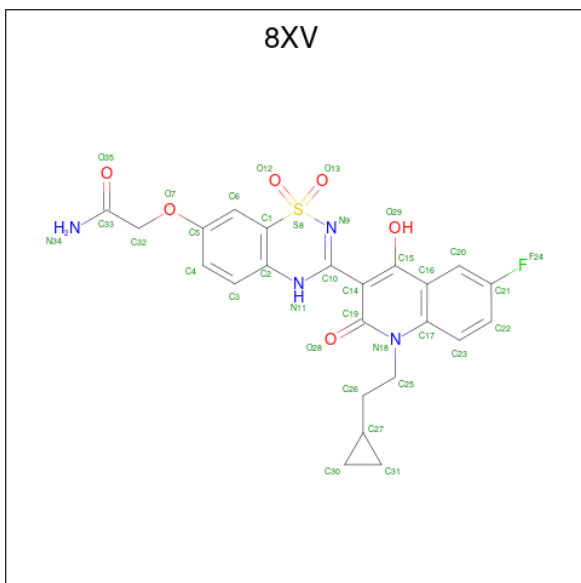
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q9WMX2
B	0	MET	-	expression tag	UNP Q9WMX2

- Molecule 2 is (2E)-3-(4-((1-((13-cyclohexyl-6-oxo-6,7-dihydro-5H-indolo[1,2-d][1,4]benzodiazepin-10-yl)carbonyl]amino)cyclopentyl)carbonyl]amino)phenyl)prop-2-enoic acid (CCD ID: 23E) (formula: C₃₈H₃₈N₄O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	38	0
			85	38	38	4	5		
2	B	1	Total	C	H	N	O	38	0
			85	38	38	4	5		

- Molecule 3 is 2-(3-[1-(2-cyclopropylethyl)-6-fluoro-4-hydroxy-2-oxo-1,2-dihydroquinolin-3-yl]-1,1-dioxo-1,4-dihydro-1λ⁶,2,4-benzothiadiazin-7-yl}oxy)acetamide (CCD ID: 8XV) (formula: C₂₃H₂₁FN₄O₆S).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	F	H	N	O	S	21	0
			56	23	1	21	4	6	1		
3	B	1	Total	C	F	H	N	O	S	21	0
			56	23	1	21	4	6	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		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

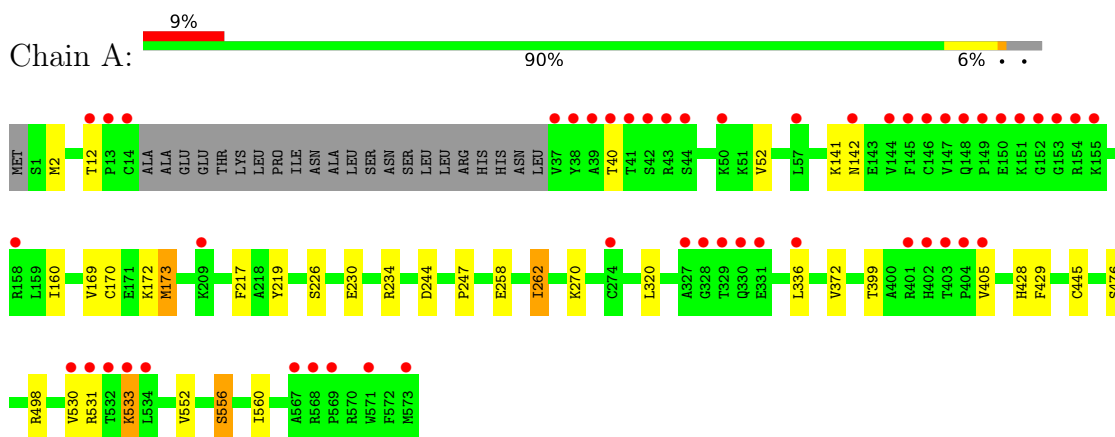
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	376	Total	O	0	0
			376	376		
5	B	377	Total	O	0	0
			377	377		

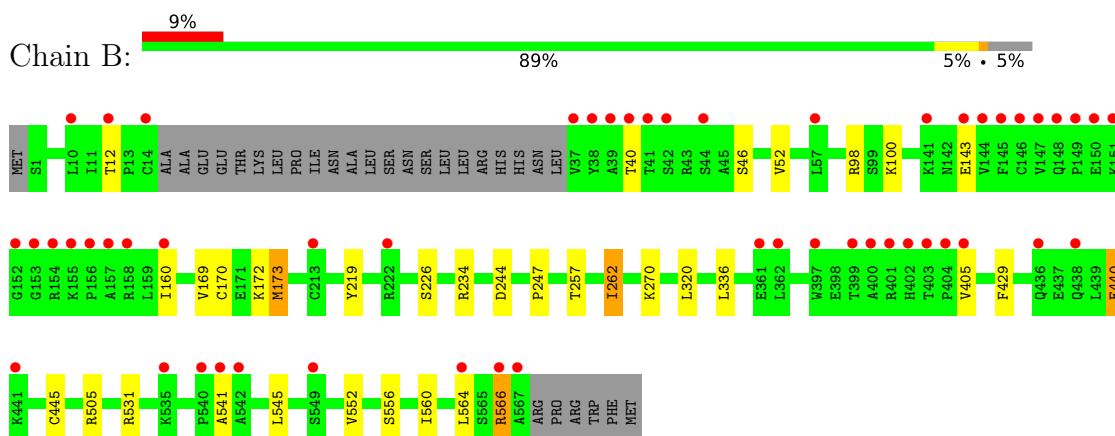
3 Residue-property plots [i](#)

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: RNA-directed RNA polymerase



• Molecule 1: RNA-directed RNA polymerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.20Å 91.30Å 232.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.44 – 2.06 41.44 – 2.06	Depositor EDS
% Data completeness (in resolution range)	96.2 (41.44-2.06) 96.1 (41.44-2.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.00 (at 2.06Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.197 , 0.228 0.198 , 0.227	Depositor DCC
R_{free} test set	1769 reflections (1.73%)	wwPDB-VP
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.633	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9492	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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: SO4, 8XV, 23E

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	0.80	2/4335 (0.0%)	1.15	5/5889 (0.1%)
1	B	0.80	2/4272 (0.0%)	1.15	3/5804 (0.1%)
All	All	0.80	4/8607 (0.0%)	1.15	8/11693 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	173	MET	SD-CE	-8.41	1.58	1.79
1	A	173	MET	SD-CE	-7.58	1.60	1.79
1	B	160	ILE	CG1-CD1	5.21	1.72	1.51
1	A	2	MET	SD-CE	5.13	1.92	1.79

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	429	PHE	CA-CB-CG	-6.58	107.22	113.80
1	A	552	VAL	N-CA-C	-6.22	106.21	111.56
1	A	429	PHE	CA-CB-CG	-6.01	107.78	113.80
1	B	552	VAL	N-CA-C	-5.68	106.68	111.56
1	A	530	VAL	CA-C-N	5.61	128.58	120.79
1	A	530	VAL	C-N-CA	5.61	128.58	120.79
1	B	244	ASP	N-CA-C	-5.18	101.32	109.25
1	A	244	ASP	N-CA-C	-5.16	101.35	109.25

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	4240	0	4185	17	0
1	B	4182	0	4137	18	0
2	A	47	38	37	0	0
2	B	47	38	37	0	0
3	A	35	21	0	0	0
3	B	35	21	0	0	0
4	A	15	0	0	0	0
4	B	20	0	0	0	0
5	A	376	0	0	1	0
5	B	377	0	0	0	0
All	All	9374	118	8396	32	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 (32) 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:170:CYS:HA	1:B:173:MET:HE3	1.59	0.84
1:A:170:CYS:HA	1:A:173:MET:HE3	1.60	0.83
1:B:545:LEU:HD22	1:B:564:LEU:HD11	1.73	0.69
1:B:505:ARG:HH12	1:B:531:ARG:HG3	1.58	0.67
1:A:141:LYS:HD2	1:A:160:ILE:HD11	1.83	0.60
1:A:247:PRO:HG3	1:B:234:ARG:HE	1.67	0.60
1:A:230:GLU:HG3	1:A:262:ILE:HG12	1.87	0.56
1:B:505:ARG:NH1	1:B:531:ARG:HG3	2.21	0.55
1:B:40:THR:HG22	1:B:143:GLU:H	1.71	0.54
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.91	0.53
1:A:12:THR:HG21	1:A:270:LYS:HG3	1.90	0.53
1:A:247:PRO:HG3	1:B:234:ARG:NE	2.23	0.52
1:B:12:THR:HG21	1:B:270:LYS:HG3	1.92	0.52
1:B:219:TYR:HB3	1:B:320:LEU:HD23	1.93	0.50
1:B:440:GLU:H	1:B:440:GLU:CD	2.21	0.48
1:B:541:ALA:HB1	1:B:566:ARG:NH2	2.31	0.46
1:A:169:VAL:HG12	1:A:173:MET:HE2	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:LYS:HE3	1:B:560:ILE:HD13	1.99	0.44
1:B:405:VAL:HG13	1:B:445:CYS:HA	1.99	0.44
1:A:556:SER:HB2	5:A:743:HOH:O	2.16	0.44
1:A:234:ARG:NH1	1:B:247:PRO:HG3	2.33	0.43
1:B:234:ARG:HG3	1:B:262:ILE:HD11	2.01	0.43
1:B:52:VAL:HB	1:B:226:SER:OG	2.19	0.42
1:A:405:VAL:HG13	1:A:445:CYS:HA	2.01	0.42
1:A:399:THR:HG21	1:A:428:HIS:NE2	2.35	0.42
1:A:234:ARG:NH1	1:A:258:GLU:OE2	2.53	0.42
1:B:169:VAL:HG12	1:B:173:MET:HE2	2.01	0.41
1:A:52:VAL:HB	1:A:226:SER:OG	2.19	0.41
1:A:172:LYS:HE3	1:A:560:ILE:HD13	2.01	0.41
1:B:257:THR:O	1:B:262:ILE:HB	2.21	0.41
1:A:40:THR:O	1:A:142:ASN:HA	2.21	0.41
1:A:217:PHE:CD2	1:A:336:LEU:HD21	2.56	0.40

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	547/574 (95%)	538 (98%)	8 (2%)	1 (0%)	43	38
1	B	541/574 (94%)	530 (98%)	11 (2%)	0	100	100
All	All	1088/1148 (95%)	1068 (98%)	19 (2%)	1 (0%)	48	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	533	LYS

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	450/486 (93%)	443 (98%)	7 (2%)	55	55
1	B	444/486 (91%)	436 (98%)	8 (2%)	51	51
All	All	894/972 (92%)	879 (98%)	15 (2%)	53	53

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

Mol	Chain	Res	Type
1	A	262	ILE
1	A	372	VAL
1	A	476	SER
1	A	498	ARG
1	A	531	ARG
1	A	533	LYS
1	A	556	SER
1	B	46	SER
1	B	98	ARG
1	B	100	LYS
1	B	262	ILE
1	B	336	LEU
1	B	440	GLU
1	B	556	SER
1	B	566	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	402	HIS
1	B	272	GLN
1	B	436	GLN

5.3.3 RNA

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](#)

11 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	$\# Z > 2$
4	SO4	A	603	-	4,4,4	0.18	0	6,6,6	0.18	0
3	8XV	B	602	-	39,39,39	0.78	1 (2%)	52,59,59	0.53	1 (1%)
4	SO4	B	606	-	4,4,4	0.25	0	6,6,6	0.08	0
4	SO4	B	604	-	4,4,4	0.25	0	6,6,6	0.05	0
2	23E	B	601	-	52,53,53	0.40	0	68,77,77	0.57	0
4	SO4	B	603	-	4,4,4	0.26	0	6,6,6	0.16	0
4	SO4	A	604	-	4,4,4	0.32	0	6,6,6	0.20	0
4	SO4	B	605	-	4,4,4	0.26	0	6,6,6	0.16	0
4	SO4	A	605	-	4,4,4	0.17	0	6,6,6	0.21	0
3	8XV	A	602	-	39,39,39	0.75	1 (2%)	52,59,59	0.55	1 (1%)
2	23E	A	601	-	52,53,53	0.37	0	68,77,77	0.51	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
2	23E	B	601	-	-	1/28/57/57	0/6/7/7
3	8XV	B	602	-	-	1/10/31/31	0/5/5/5
3	8XV	A	602	-	-	1/10/31/31	0/5/5/5
2	23E	A	601	-	-	0/28/57/57	0/6/7/7

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	8XV	C14-C10	3.24	1.52	1.47
3	A	602	8XV	C14-C10	2.99	1.52	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	8XV	C25-C26-C27	-2.20	109.98	114.85
3	A	602	8XV	C25-C26-C27	-2.18	110.03	114.85

There are no chirality outliers.

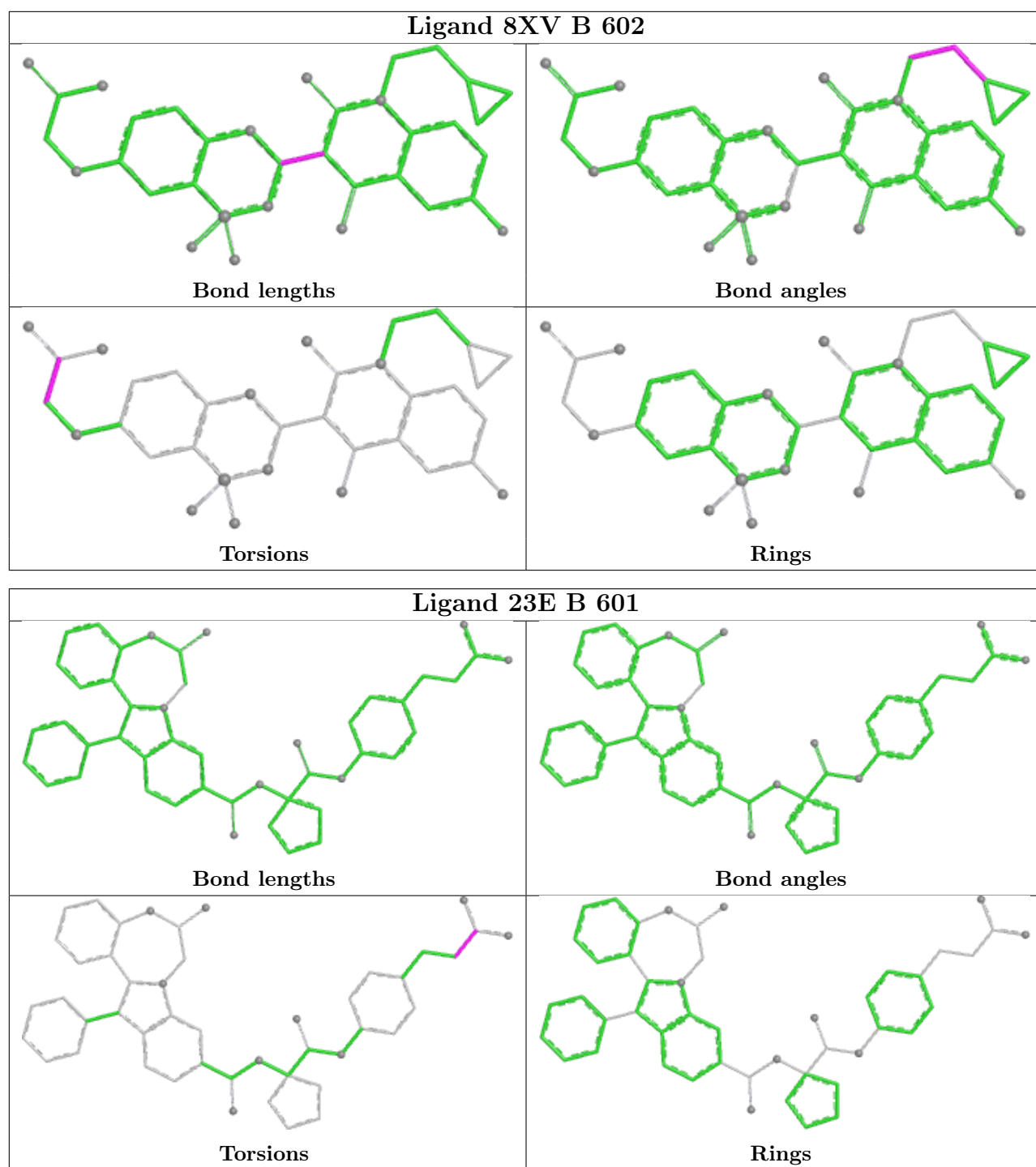
All (3) torsion outliers are listed below:

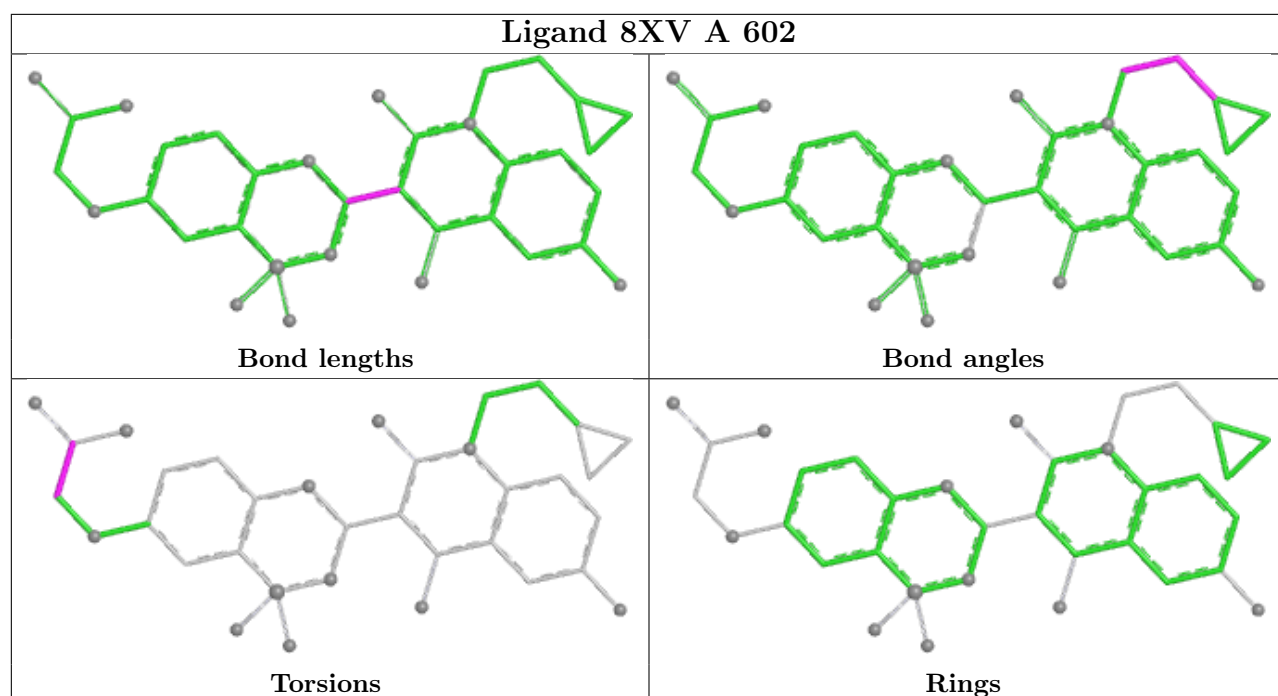
Mol	Chain	Res	Type	Atoms
3	A	602	8XV	O7-C32-C33-O35
3	B	602	8XV	O7-C32-C33-O35
2	B	601	23E	C12-C13-C14-O15

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	551/574 (95%)	0.33	50 (9%)	15 15	12, 23, 59, 101	0
1	B	545/574 (94%)	0.51	52 (9%)	14 14	13, 26, 64, 107	0
All	All	1096/1148 (95%)	0.42	102 (9%)	14 14	12, 24, 60, 107	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	567	ALA	7.7
1	A	149	PRO	6.5
1	A	147	VAL	6.4
1	B	405	VAL	6.2
1	B	37	VAL	6.1
1	B	149	PRO	5.7
1	B	146	CYS	5.6
1	A	37	VAL	5.4
1	B	148	GLN	5.4
1	A	14	CYS	5.2
1	B	145	PHE	5.0
1	B	38	TYR	5.0
1	A	148	GLN	5.0
1	B	42	SER	5.0
1	B	403	THR	4.9
1	B	151	LYS	4.9
1	B	40	THR	4.9
1	B	147	VAL	4.9
1	B	404	PRO	4.8
1	A	150	GLU	4.7
1	A	532	THR	4.7
1	B	150	GLU	4.5
1	B	157	ALA	4.4
1	B	155	LYS	4.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	41	THR	4.3
1	B	564	LEU	4.3
1	B	14	CYS	4.3
1	B	400	ALA	4.3
1	B	402	HIS	4.1
1	B	154	ARG	3.9
1	A	39	ALA	3.9
1	B	41	THR	3.9
1	A	151	LYS	3.9
1	B	535	LYS	3.9
1	B	39	ALA	3.9
1	A	573	MET	3.9
1	B	153	GLY	3.8
1	A	533	LYS	3.8
1	A	327	ALA	3.6
1	A	403	THR	3.5
1	A	571	TRP	3.4
1	B	401	ARG	3.4
1	A	569	PRO	3.4
1	A	153	GLY	3.3
1	B	152	GLY	3.3
1	A	146	CYS	3.2
1	A	145	PHE	3.2
1	A	404	PRO	3.2
1	B	156	PRO	3.2
1	B	144	VAL	3.1
1	A	40	THR	3.1
1	B	541	ALA	3.1
1	B	44	SER	3.0
1	A	142	ASN	3.0
1	A	568	ARG	2.9
1	A	274	CYS	2.9
1	B	57	LEU	2.9
1	A	567	ALA	2.9
1	A	152	GLY	2.9
1	B	542	ALA	2.9
1	A	13	PRO	2.8
1	A	38	TYR	2.8
1	A	531	ARG	2.8
1	B	397	TRP	2.8
1	A	154	ARG	2.8
1	A	12	THR	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	160	ILE	2.8
1	A	330	GLN	2.8
1	B	143	GLU	2.7
1	A	534	LEU	2.7
1	A	144	VAL	2.7
1	A	530	VAL	2.7
1	B	549	SER	2.7
1	A	336	LEU	2.7
1	B	566	ARG	2.7
1	B	141	LYS	2.6
1	B	158	ARG	2.6
1	A	42	SER	2.6
1	A	43	ARG	2.5
1	A	405	VAL	2.5
1	B	399	THR	2.5
1	A	155	LYS	2.5
1	A	402	HIS	2.4
1	A	209	LYS	2.4
1	A	401	ARG	2.4
1	B	540	PRO	2.4
1	A	328	GLY	2.3
1	A	44	SER	2.2
1	B	436	GLN	2.2
1	A	329	THR	2.2
1	A	331	GLU	2.1
1	B	361	GLU	2.1
1	A	158	ARG	2.1
1	B	213	CYS	2.1
1	B	12	THR	2.1
1	A	50	LYS	2.1
1	B	10	LEU	2.0
1	B	362	LEU	2.0
1	B	438	GLN	2.0
1	A	57	LEU	2.0
1	B	222	ARG	2.0
1	B	441	LYS	2.0

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

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](#)

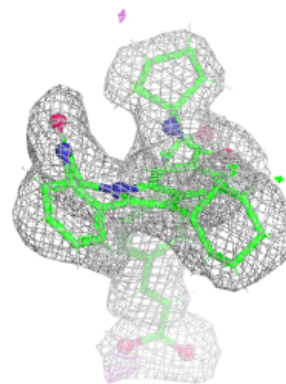
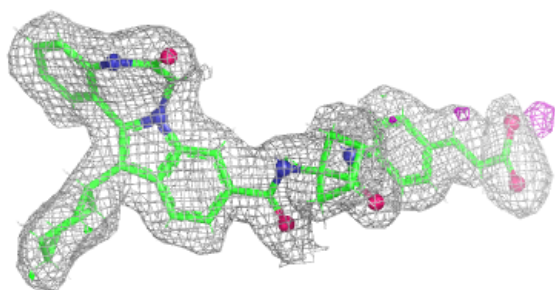
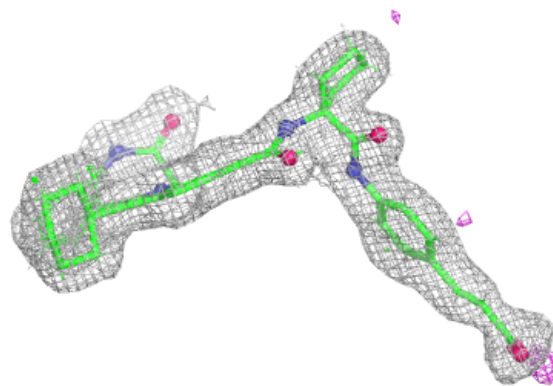
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	SO4	B	603	5/5	0.71	0.16	97,98,98,98	0
4	SO4	A	604	5/5	0.74	0.17	75,77,78,79	0
4	SO4	B	604	5/5	0.78	0.15	101,101,102,102	0
4	SO4	B	605	5/5	0.79	0.20	112,112,114,114	0
4	SO4	B	606	5/5	0.83	0.14	87,88,88,89	0
4	SO4	A	603	5/5	0.86	0.10	63,64,64,66	0
4	SO4	A	605	5/5	0.90	0.12	56,57,59,61	0
2	23E	B	601	47/47	0.92	0.09	15,21,35,42	38
2	23E	A	601	47/47	0.94	0.08	13,17,41,46	38
3	8XV	B	602	35/35	0.97	0.06	16,19,32,33	21
3	8XV	A	602	35/35	0.97	0.06	12,18,23,26	21

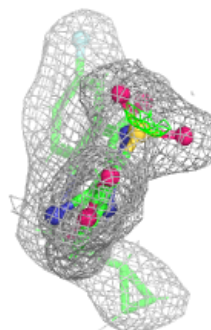
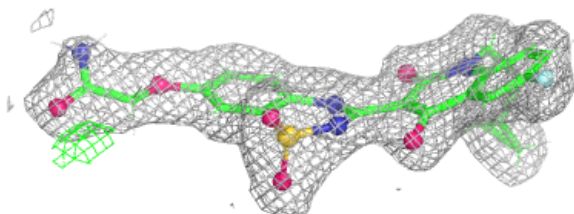
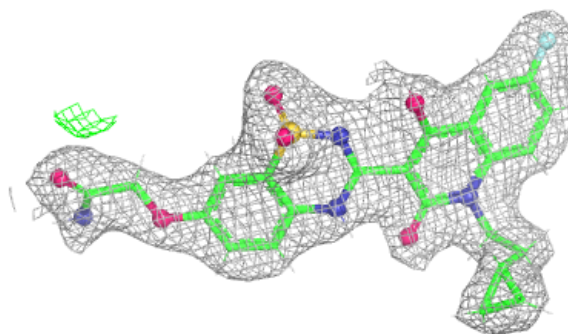
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 23E B 601:

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

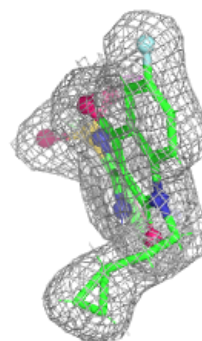
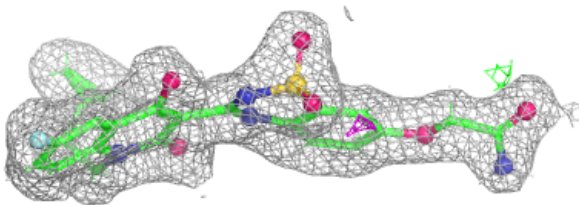
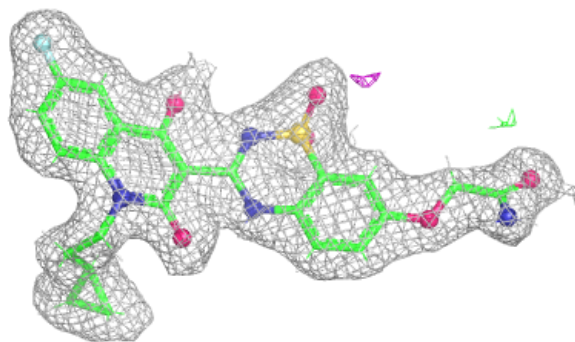
**Electron density around 8XV B 602:**

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



Electron density around 8XV A 602:

$2mF_o-DF_c$ (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.