



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:43 PM UTC

PDB ID : 2AWZ / pdb_00002awz
Title : Hepatitis C Virus NS5b RNA Polymerase in complex with a covalent inhibitor (5h)
Authors : Powers, J.P.; Piper, D.E.; Li, Y.; Mayorga, V.; Anzola, J.; Chen, J.M.; Jaen, J.C.; Lee, G.; Liu, J.; Peterson, M.G.; Tonn, G.R.; Ye, Q.; Walker, N.P.; Wang, Z.
Deposited on : 2005-09-02
Resolution : 2.15 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](#) ①) 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

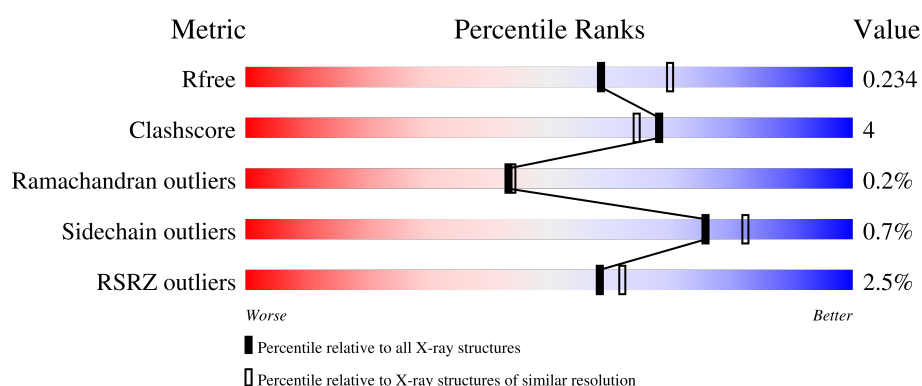
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	580	2% 85% 11% .
1	B	580	3% 88% 9% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9525 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 Genome polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4337	2735	767	804	31			
1	B	562	Total	C	N	O	S	0	0	0
			4370	2755	775	809	31			

There are 42 discrepancies between the modelled and reference sequences:

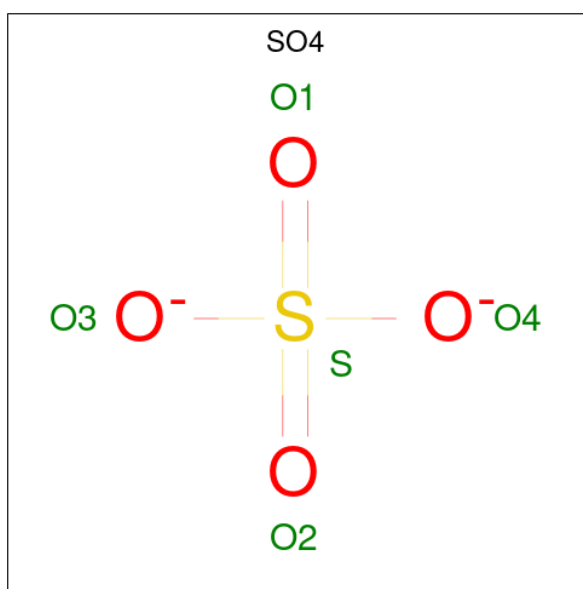
Chain	Residue	Modelled	Actual	Comment	Reference
A	499	ALA	VAL	engineered mutation	UNP P26663
A	506	ASN	SER	engineered mutation	UNP P26663
A	514	ARG	GLN	engineered mutation	UNP P26663
A	520	ILE	THR	engineered mutation	UNP P26663
A	540	ALA	PRO	engineered mutation	UNP P26663
A	543	GLY	SER	engineered mutation	UNP P26663
A	549	SER	GLY	engineered mutation	UNP P26663
A	552	THR	VAL	engineered mutation	UNP P26663
A	563	GLY	SER	engineered mutation	UNP P26663
A	564	VAL	LEU	engineered mutation	UNP P26663
A	566	HIS	ARG	engineered mutation	UNP P26663
A	571	HIS	-	expression tag	UNP P26663
A	572	HIS	-	expression tag	UNP P26663
A	573	HIS	-	expression tag	UNP P26663
A	574	HIS	-	expression tag	UNP P26663
A	575	HIS	-	expression tag	UNP P26663
A	576	HIS	-	expression tag	UNP P26663
A	577	HIS	-	expression tag	UNP P26663
A	578	HIS	-	expression tag	UNP P26663
A	579	HIS	-	expression tag	UNP P26663
A	580	HIS	-	expression tag	UNP P26663
B	499	ALA	VAL	engineered mutation	UNP P26663
B	506	ASN	SER	engineered mutation	UNP P26663
B	514	ARG	GLN	engineered mutation	UNP P26663
B	520	ILE	THR	engineered mutation	UNP P26663

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Chain	Residue	Modelled	Actual	Comment	Reference
B	540	ALA	PRO	engineered mutation	UNP P26663
B	543	GLY	SER	engineered mutation	UNP P26663
B	549	SER	GLY	engineered mutation	UNP P26663
B	552	THR	VAL	engineered mutation	UNP P26663
B	563	GLY	SER	engineered mutation	UNP P26663
B	564	VAL	LEU	engineered mutation	UNP P26663
B	566	HIS	ARG	engineered mutation	UNP P26663
B	571	HIS	-	expression tag	UNP P26663
B	572	HIS	-	expression tag	UNP P26663
B	573	HIS	-	expression tag	UNP P26663
B	574	HIS	-	expression tag	UNP P26663
B	575	HIS	-	expression tag	UNP P26663
B	576	HIS	-	expression tag	UNP P26663
B	577	HIS	-	expression tag	UNP P26663
B	578	HIS	-	expression tag	UNP P26663
B	579	HIS	-	expression tag	UNP P26663
B	580	HIS	-	expression tag	UNP P26663

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



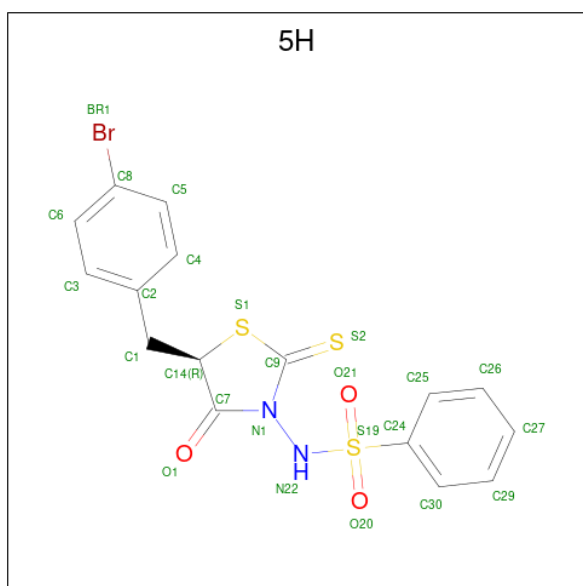
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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

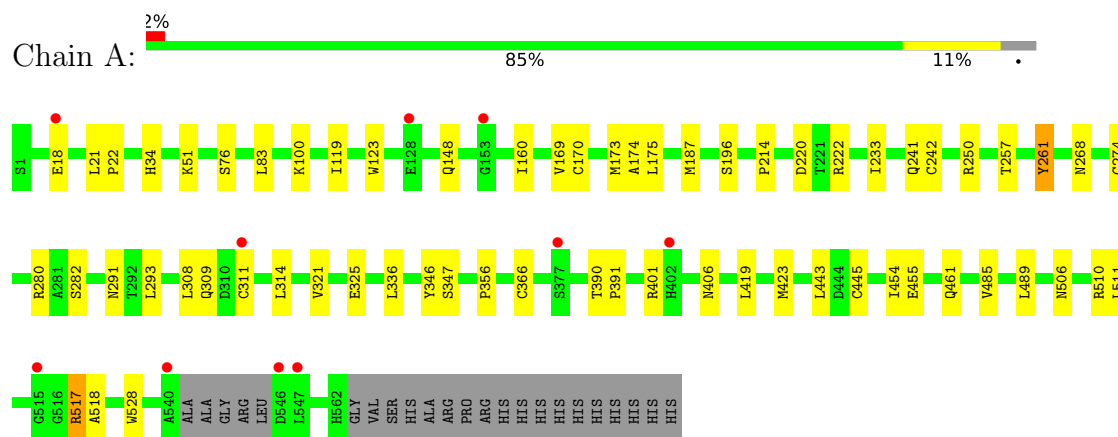
- Molecule 3 is 5R-(4-BROMOPHENYLMETHYL)-3-(BENZENESULFONYLAMINO)-4-OXO-2-THIONOTHIAZOLIDINE (CCD ID: 5H) (formula: C₁₆H₁₃BrN₂O₃S₃).



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: Genome polyprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.09Å 87.02Å 162.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.15 25.00 – 2.16	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.15) 96.4 (25.00-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.15Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.205 , 0.243 0.197 , 0.234	Depositor DCC
R_{free} test set	6552 reflections (9.71%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9525	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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, 5H

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	0.40	0/4431	0.85	6/6010 (0.1%)
1	B	0.40	0/4465	0.85	4/6057 (0.1%)
All	All	0.40	0/8896	0.85	10/12067 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	SER	N-CA-C	-9.31	99.43	110.13
1	A	261	TYR	N-CA-C	6.92	118.82	111.28
1	B	196	SER	N-CA-C	-6.78	100.17	110.07
1	B	151	LYS	N-CA-C	6.05	123.69	110.80
1	A	175	LEU	N-CA-C	5.73	122.69	114.39
1	B	261	TYR	N-CA-C	5.51	117.29	111.28
1	A	148	GLN	CA-C-N	5.42	124.88	119.24
1	A	148	GLN	C-N-CA	5.42	124.88	119.24
1	B	174	ALA	N-CA-C	5.16	118.04	111.69
1	A	214	PRO	N-CA-C	5.11	119.11	111.14

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	4337	0	4355	35	0
1	B	4370	0	4393	28	0
2	A	20	0	0	0	0
2	B	10	0	0	0	0
3	A	25	0	12	0	0
3	B	25	0	12	0	0
4	A	364	0	0	2	0
4	B	374	0	0	0	0
All	All	9525	0	8772	63	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 (63) 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:517:ARG:HH11	1:A:517:ARG:HG3	1.40	0.86
1:A:83:LEU:HB2	1:A:173:MET:HA	1.68	0.74
1:B:233:ILE:HD13	1:B:262:ILE:HA	1.69	0.73
1:B:83:LEU:HB2	1:B:173:MET:HA	1.72	0.70
1:B:336:LEU:HD22	1:B:356:PRO:HG3	1.75	0.69
1:A:241:GLN:OE1	1:A:250:ARG:HG3	1.92	0.69
1:A:445:CYS:SG	1:A:454:ILE:HD12	2.33	0.69
1:A:18:GLU:HG2	1:A:401:ARG:CZ	2.25	0.66
1:B:150:GLU:O	1:B:152:GLY:N	2.29	0.64
1:A:506:ASN:HD21	1:A:510:ARG:HH11	1.49	0.61
1:A:336:LEU:HD22	1:A:356:PRO:HG3	1.84	0.59
1:B:313:MET:HG2	1:B:322:VAL:HG22	1.84	0.59
1:A:455:GLU:HB3	4:A:1033:HOH:O	2.03	0.59
1:A:280:ARG:HD2	1:A:291:ASN:OD1	2.03	0.58
1:A:308:LEU:HB2	1:A:311:CYS:SG	2.46	0.55
1:A:517:ARG:HG3	1:A:517:ARG:NH1	2.16	0.55
1:A:119:ILE:HD13	1:A:169:VAL:HG11	1.89	0.54
1:B:172:LYS:HE3	1:B:560:ILE:HD13	1.90	0.53
1:B:346:TYR:O	1:B:347:SER:HB3	2.07	0.53
1:B:361:GLU:OE1	1:B:372:VAL:HG23	2.09	0.53
1:B:309:GLN:HG3	1:B:325:GLU:HB2	1.91	0.52
1:A:34:HIS:HB3	4:A:1244:HOH:O	2.10	0.52
1:B:308:LEU:HB2	1:B:311:CYS:SG	2.50	0.51
1:B:115:ALA:O	1:B:119:ILE:HG13	2.11	0.51
1:B:82:LEU:HD13	1:B:249:ALA:HB2	1.93	0.51
1:A:346:TYR:O	1:A:347:SER:HB3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:ARG:NH2	1:B:279:CYS:HB3	2.26	0.50
1:B:183:PRO:HG3	1:B:289:CYS:SG	2.52	0.49
1:A:419:LEU:HD11	1:A:485:VAL:HG11	1.94	0.49
1:A:51:LYS:HG3	1:A:222:ARG:NH2	2.28	0.48
1:A:18:GLU:HG2	1:A:401:ARG:NH2	2.29	0.48
1:B:237:GLU:OE2	1:B:254:LYS:HA	2.15	0.47
1:A:233:ILE:HD13	1:A:261:TYR:O	2.15	0.47
1:A:160:ILE:HA	1:A:282:SER:OG	2.16	0.46
1:A:308:LEU:CB	1:A:311:CYS:SG	3.03	0.46
1:B:187:MET:HE3	1:B:296:TYR:CD1	2.50	0.46
1:A:268:ASN:HB3	1:A:274:CYS:SG	2.56	0.46
1:B:515:GLY:HA2	1:B:519:ALA:HB2	1.98	0.45
1:A:390:THR:HB	1:A:391:PRO:HD3	1.98	0.45
1:A:485:VAL:O	1:A:489:LEU:HG	2.17	0.45
1:A:187:MET:HE2	1:A:293:LEU:HD23	1.98	0.44
1:B:390:THR:HB	1:B:391:PRO:HD3	1.99	0.44
1:B:355:GLN:HA	1:B:356:PRO:HD3	1.90	0.44
1:B:170:CYS:HA	1:B:173:MET:HE3	2.00	0.44
1:B:187:MET:HE1	1:B:292:THR:HG22	1.99	0.44
1:B:334:ALA:HA	1:B:337:ARG:HH11	1.83	0.43
1:B:429:PHE:O	1:B:433:LEU:HG	2.18	0.43
1:A:257:THR:HA	1:A:261:TYR:HB2	2.01	0.43
1:A:21:LEU:HD12	1:A:22:PRO:HD2	2.02	0.42
1:A:76:SER:HA	1:A:242:CYS:O	2.18	0.42
1:A:511:LEU:HB3	1:A:518:ALA:O	2.19	0.42
1:A:100:LYS:HE2	1:A:100:LYS:HB3	1.86	0.42
1:B:149:PRO:O	1:B:150:GLU:O	2.38	0.42
1:B:309:GLN:H	1:B:309:GLN:HG2	1.73	0.41
1:A:406:ASN:ND2	1:A:443:LEU:HB3	2.35	0.41
1:A:423:MET:HA	1:A:528:TRP:CZ2	2.55	0.41
1:B:423:MET:HA	1:B:528:TRP:CZ2	2.55	0.41
1:B:56:ARG:NH2	1:B:228:VAL:O	2.54	0.41
1:A:309:GLN:HB2	1:A:325:GLU:HB2	2.02	0.41
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.51	0.40
1:B:216:GLY:HA2	1:B:322:VAL:O	2.21	0.40
1:A:123:TRP:CH2	1:A:174:ALA:HB2	2.57	0.40
1:A:170:CYS:HA	1:A:173:MET:HE3	2.04	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	553/580 (95%)	544 (98%)	9 (2%)	0	100	100
1	B	560/580 (97%)	551 (98%)	7 (1%)	2 (0%)	30	26
All	All	1113/1160 (96%)	1095 (98%)	16 (1%)	2 (0%)	43	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	151	LYS
1	B	150	GLU

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	474/492 (96%)	470 (99%)	4 (1%)	73	79
1	B	476/492 (97%)	473 (99%)	3 (1%)	78	84
All	All	950/984 (96%)	943 (99%)	7 (1%)	76	82

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	220	ASP
1	A	366	CYS
1	A	461	GLN
1	A	517	ARG

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Mol	Chain	Res	Type
1	B	56	ARG
1	B	150	GLU
1	B	366	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	35	ASN
1	A	95	HIS
1	A	184	GLN
1	A	206	ASN
1	A	213	ASN
1	A	309	GLN
1	A	406	ASN
1	A	438	GLN
1	A	461	GLN
1	A	506	ASN
1	B	330	GLN
1	B	506	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](#)

8 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	SO4	A	903	-	4,4,4	0.43	0	6,6,6	0.09	0
2	SO4	B	905	-	4,4,4	0.45	0	6,6,6	0.11	0
3	5H	B	802	1	25,27,27	1.80	3 (12%)	32,39,39	1.18	2 (6%)
2	SO4	A	902	-	4,4,4	0.36	0	6,6,6	0.05	0
2	SO4	B	906	-	4,4,4	0.46	0	6,6,6	0.08	0
2	SO4	A	904	-	4,4,4	0.45	0	6,6,6	0.08	0
2	SO4	A	901	-	4,4,4	0.43	0	6,6,6	0.06	0
3	5H	A	801	1	25,27,27	1.72	3 (12%)	32,39,39	1.35	3 (9%)

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	5H	A	801	1	-	2/13/31/31	0/3/3/3
3	5H	B	802	1	-	1/13/31/31	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	5H	S19-N22	4.66	1.75	1.65
3	A	801	5H	S19-N22	4.44	1.74	1.65
3	B	802	5H	C14-S1	-3.76	1.75	1.82
3	A	801	5H	C7-N1	-3.34	1.33	1.39
3	B	802	5H	C7-N1	-3.32	1.33	1.39
3	A	801	5H	C14-S1	-2.79	1.77	1.82

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	5H	C7-C14-S1	-4.68	102.38	106.11
3	A	801	5H	C7-N1-N22	4.18	126.63	122.74
3	B	802	5H	C7-N1-N22	3.57	126.06	122.74
3	B	802	5H	C7-C14-S1	-3.48	103.34	106.11
3	A	801	5H	O1-C7-N1	-2.03	121.30	123.60

There are no chirality outliers.

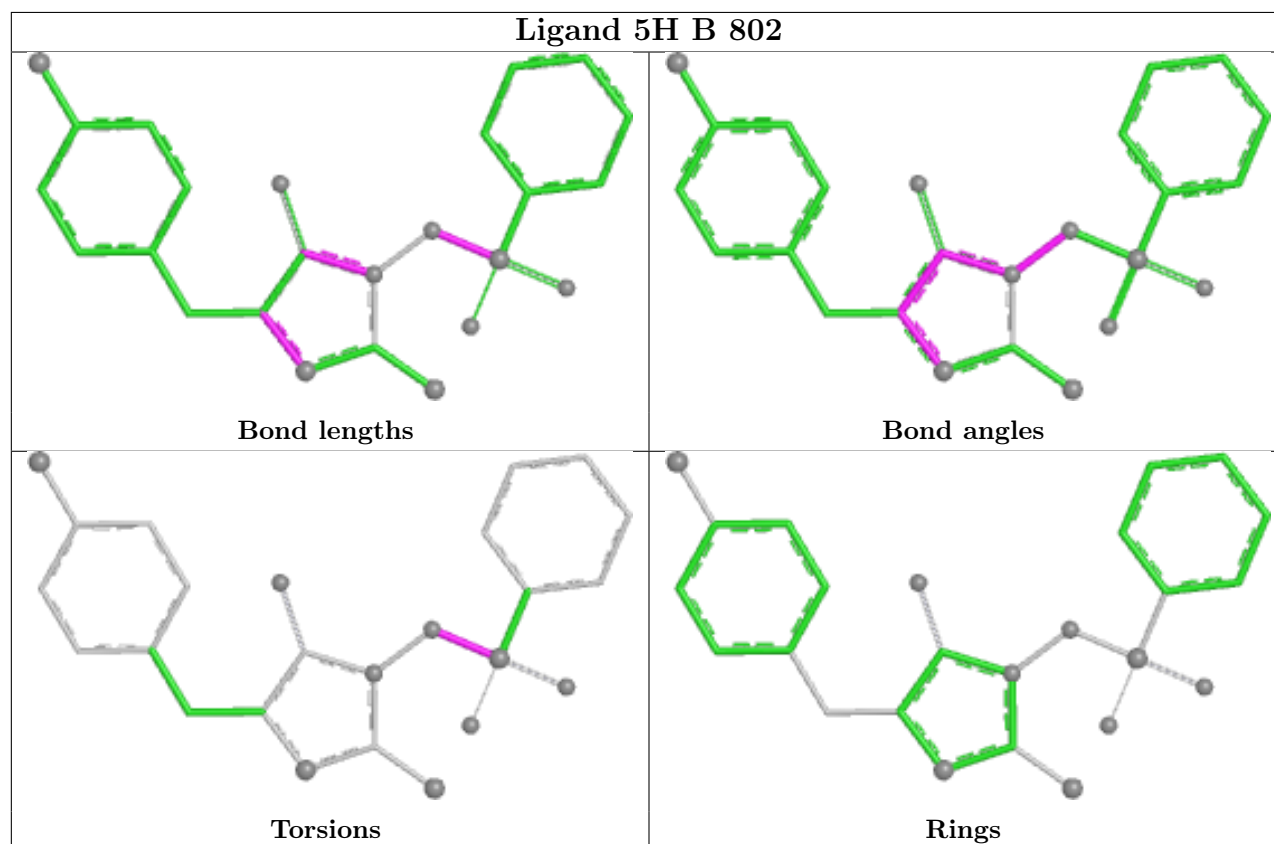
All (3) torsion outliers are listed below:

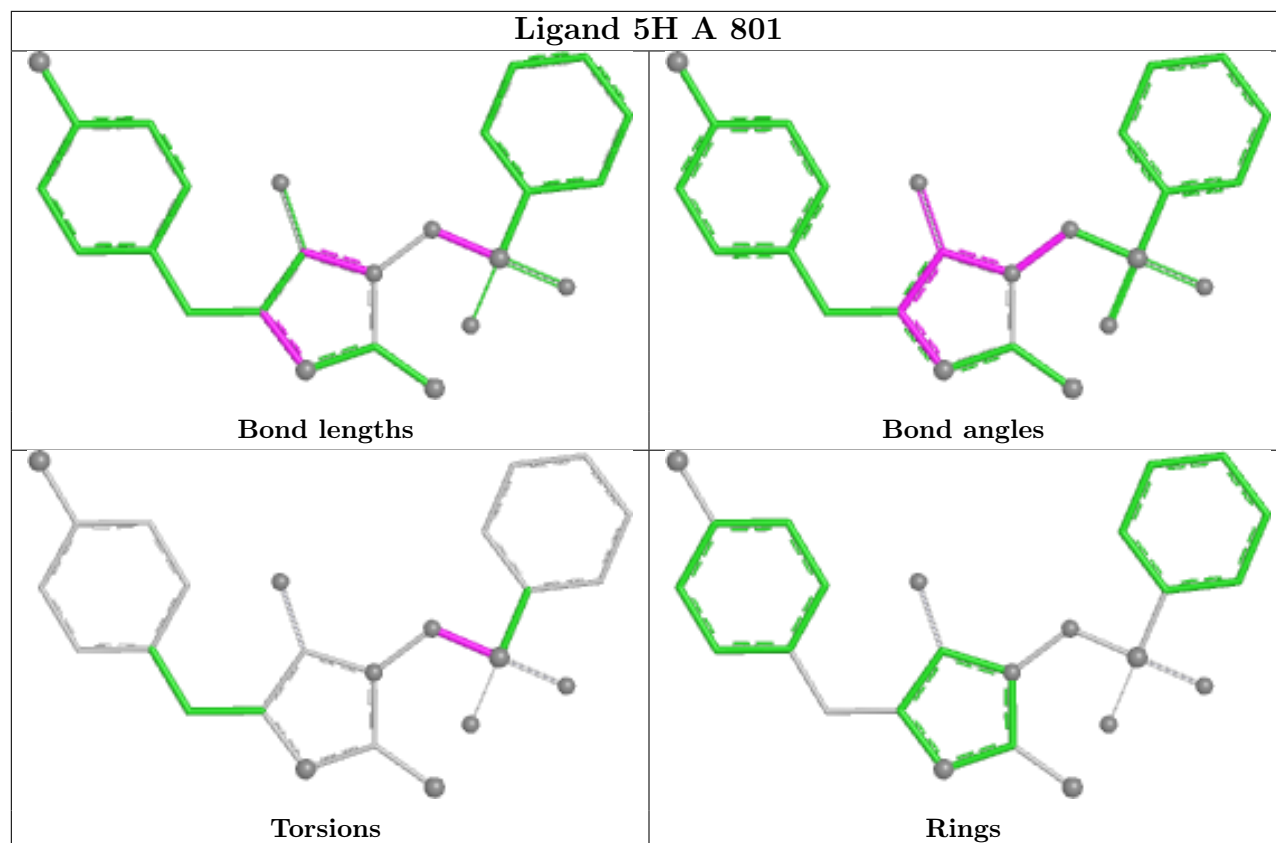
Mol	Chain	Res	Type	Atoms
3	B	802	5H	N1-N22-S19-O20
3	A	801	5H	N1-N22-S19-O20
3	A	801	5H	N1-N22-S19-O21

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	557/580 (96%)	0.10	10 (1%) 67 71	14, 25, 44, 64	0
1	B	562/580 (96%)	0.14	18 (3%) 50 54	12, 26, 45, 57	0
All	All	1119/1160 (96%)	0.12	28 (2%) 58 62	12, 26, 44, 64	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	149	PRO	3.6
1	A	547	LEU	3.5
1	B	376	ALA	3.2
1	A	515	GLY	3.1
1	B	403	THR	2.9
1	B	150	GLU	2.9
1	B	405	VAL	2.9
1	B	14	CYS	2.8
1	A	546	ASP	2.7
1	B	540	ALA	2.7
1	B	151	LYS	2.7
1	A	402	HIS	2.6
1	A	128	GLU	2.5
1	B	542	ALA	2.5
1	B	534	LEU	2.4
1	B	402	HIS	2.4
1	B	82	LEU	2.4
1	B	532	THR	2.4
1	A	540	ALA	2.3
1	A	311	CYS	2.3
1	B	406	ASN	2.3
1	B	246	ALA	2.3
1	A	18	GLU	2.2
1	B	531	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	153	GLY	2.1
1	B	90	LYS	2.1
1	A	377	SER	2.1
1	B	544	ARG	2.0

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](#)

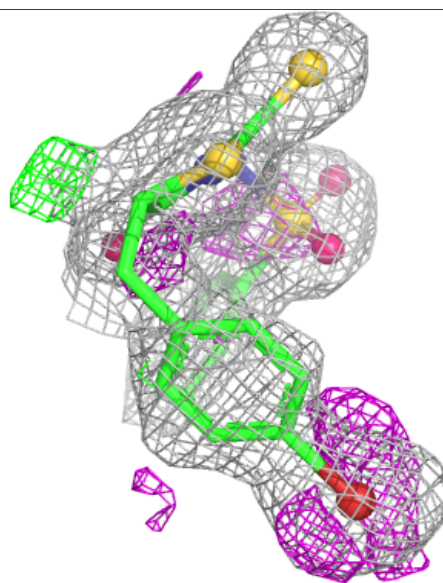
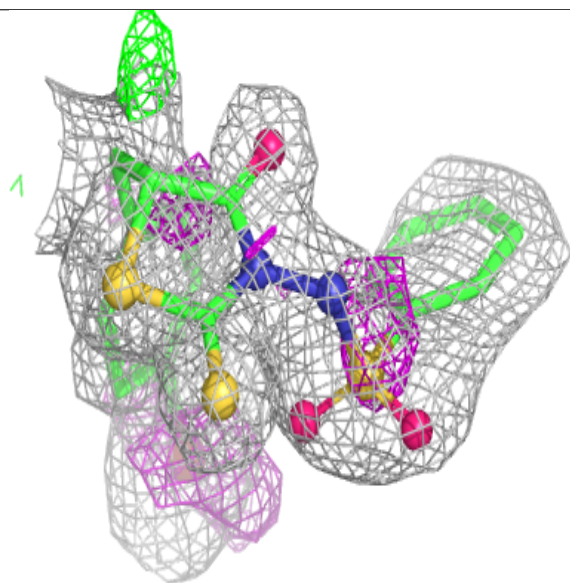
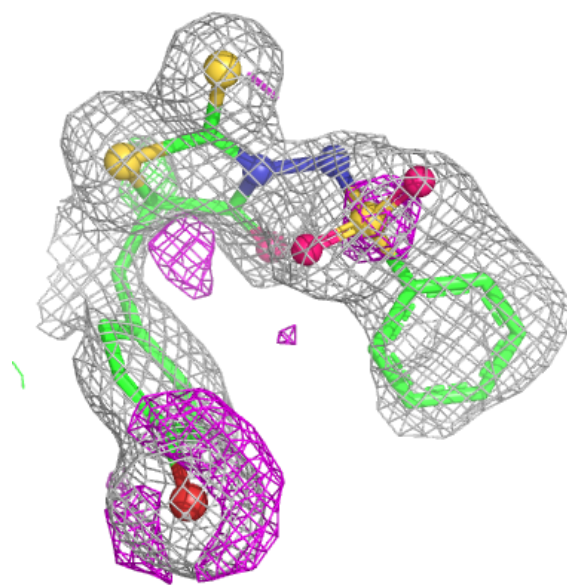
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	SO4	B	906	5/5	0.75	0.12	73,73,73,73	0
2	SO4	A	904	5/5	0.84	0.14	41,42,43,43	0
2	SO4	A	903	5/5	0.87	0.13	48,48,49,49	0
3	5H	A	801	25/25	0.92	0.12	46,51,52,56	0
2	SO4	A	902	5/5	0.93	0.08	40,40,41,42	0
3	5H	B	802	25/25	0.93	0.12	47,49,51,55	0
2	SO4	B	905	5/5	0.95	0.12	32,33,33,34	0
2	SO4	A	901	5/5	0.97	0.07	30,30,31,31	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.

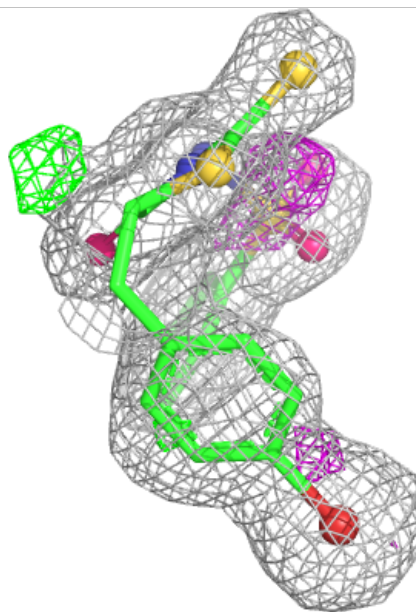
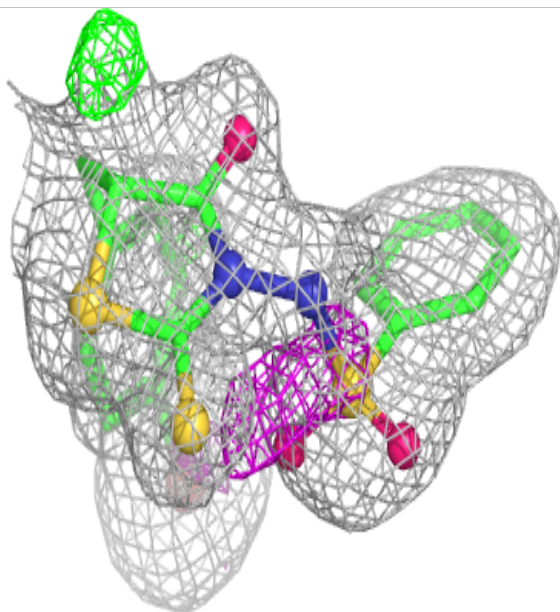
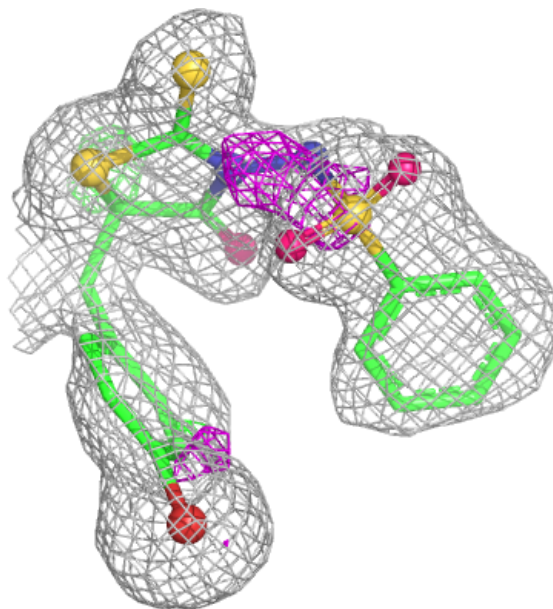
Electron density around 5H A 801:

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



Electron density around 5H B 802:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.