



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:00 pm BST

PDB ID : 4OJQ
Title : Crystal Structure of Hepatitis C Virus NS3 Helicase Inhibitor Co-complex with
Fragment 1 [(5-bromo-1H-indol-3-yl)acetic acid]
Authors : Padyana, A.K.
Deposited on : 2014-01-21
Resolution : 2.25 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

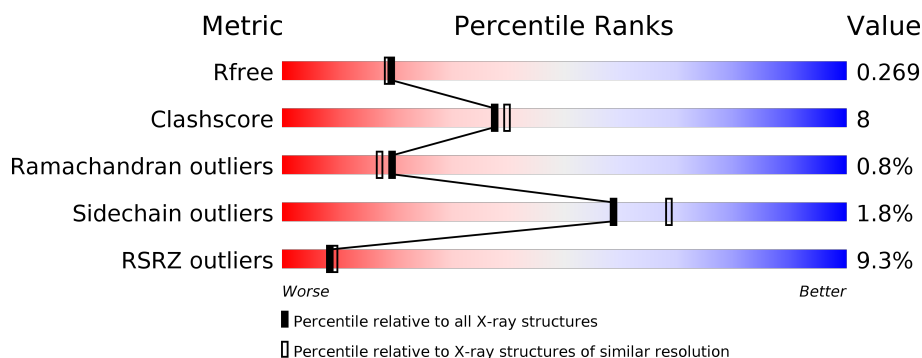
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 on 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	464	<div> <div>10%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>• 5%</div> </div> </div>
1	B	464	<div> <div>7%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>• 9%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6714 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 Serine protease NS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	4	0
			3337	2115	560	637	25			
1	B	420	Total	C	N	O	S	0	1	0
			3176	2018	532	603	23			

There are 30 discrepancies between the modelled and reference sequences:

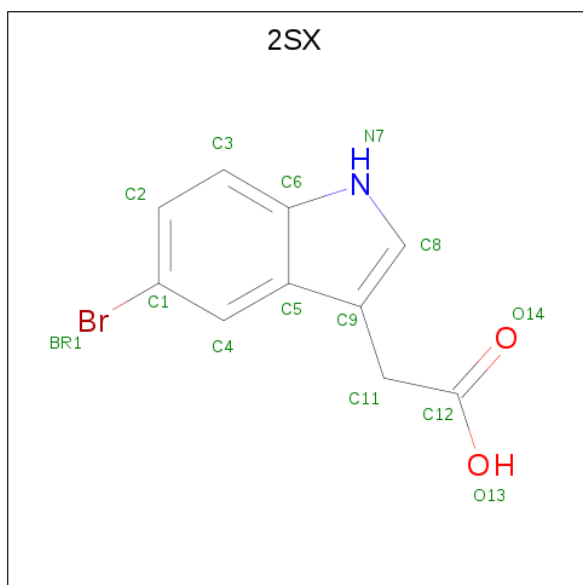
Chain	Residue	Modelled	Actual	Comment	Reference
A	167	MET	-	EXPRESSION TAG	UNP K4KA16
A	168	GLY	-	EXPRESSION TAG	UNP K4KA16
A	169	SER	-	EXPRESSION TAG	UNP K4KA16
A	170	SER	-	EXPRESSION TAG	UNP K4KA16
A	171	HIS	-	EXPRESSION TAG	UNP K4KA16
A	172	HIS	-	EXPRESSION TAG	UNP K4KA16
A	173	HIS	-	EXPRESSION TAG	UNP K4KA16
A	174	HIS	-	EXPRESSION TAG	UNP K4KA16
A	175	HIS	-	EXPRESSION TAG	UNP K4KA16
A	176	HIS	-	EXPRESSION TAG	UNP K4KA16
A	177	SER	-	EXPRESSION TAG	UNP K4KA16
A	178	SER	-	EXPRESSION TAG	UNP K4KA16
A	179	GLY	-	EXPRESSION TAG	UNP K4KA16
A	403	ASN	SER	CONFLICT	UNP K4KA16
A	505	MET	THR	CONFLICT	UNP K4KA16
B	167	MET	-	EXPRESSION TAG	UNP K4KA16
B	168	GLY	-	EXPRESSION TAG	UNP K4KA16
B	169	SER	-	EXPRESSION TAG	UNP K4KA16
B	170	SER	-	EXPRESSION TAG	UNP K4KA16
B	171	HIS	-	EXPRESSION TAG	UNP K4KA16
B	172	HIS	-	EXPRESSION TAG	UNP K4KA16
B	173	HIS	-	EXPRESSION TAG	UNP K4KA16
B	174	HIS	-	EXPRESSION TAG	UNP K4KA16
B	175	HIS	-	EXPRESSION TAG	UNP K4KA16
B	176	HIS	-	EXPRESSION TAG	UNP K4KA16

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Chain	Residue	Modelled	Actual	Comment	Reference
B	177	SER	-	EXPRESSION TAG	UNP K4KA16
B	178	SER	-	EXPRESSION TAG	UNP K4KA16
B	179	GLY	-	EXPRESSION TAG	UNP K4KA16
B	403	ASN	SER	CONFLICT	UNP K4KA16
B	505	MET	THR	CONFLICT	UNP K4KA16

- Molecule 2 is (5-bromo-1H-indol-3-yl)acetic acid (three-letter code: 2SX) (formula: $C_{10}H_8BrNO_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	0	0
			14	1	10	1	2		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Ca	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	85	Total	O	0	0
			85	85		
4	B	100	Total	O	0	0
			100	100		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.14Å 103.53Å 119.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.10 – 2.25 39.10 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.2 (39.10-2.25) 96.9 (39.10-2.25)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.217 , 0.270 0.219 , 0.269	Depositor DCC
R_{free} test set	2458 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6714	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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: CA, 2SX

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	0.45	0/3416	0.64	3/4666 (0.1%)
1	B	0.46	0/3253	0.62	1/4441 (0.0%)
All	All	0.46	0/6669	0.63	4/9107 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	588	LEU	CA-CB-CG	8.26	134.31	115.30
1	A	588	LEU	CB-CG-CD2	6.59	122.20	111.00
1	B	588	LEU	CA-CB-CG	5.75	128.54	115.30
1	A	592	LEU	CA-CB-CG	5.65	128.29	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	397	VAL	Peptide
1	A	399	VAL	Peptide

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Mol	Chain	Res	Type	Group
1	B	401	PRO	Peptide

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	3337	0	3292	56	1
1	B	3176	0	3127	46	1
2	A	14	0	7	0	0
3	B	2	0	0	0	0
4	A	85	0	0	1	0
4	B	100	0	0	3	0
All	All	6714	0	6426	100	1

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 8.

All (100) 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:363:ARG:HG3	1:A:422:PHE:HA	1.55	0.87
1:B:402:THR:HG22	1:B:403:ASN:H	1.42	0.84
1:A:397:VAL:HG21	1:A:418:PHE:HE2	1.46	0.81
1:A:399:VAL:HG12	1:A:400:ILE:HB	1.68	0.75
1:B:417:GLY:N	4:B:1199:HOH:O	2.20	0.74
1:B:460:GLN:HE21	1:B:464:ARG:HH11	1.36	0.72
1:A:397:VAL:HG21	1:A:418:PHE:CE2	2.22	0.72
1:B:402:THR:CG2	1:B:403:ASN:H	2.03	0.72
1:A:381:LEU:HA	1:A:384:LEU:HB2	1.72	0.71
1:A:392:TYR:O	1:A:395:LEU:N	2.25	0.69
1:A:396:ASP:OD1	1:A:397:VAL:N	2.22	0.69
1:A:364:HIS:HB2	1:A:407:VAL:HG22	1.80	0.64
1:A:363:ARG:CZ	1:A:400:ILE:HD11	2.29	0.63
1:A:366:ILE:HG12	1:A:426:ILE:HB	1.81	0.62
1:B:389:VAL:HG21	1:B:400:ILE:HD13	1.81	0.61
1:A:411:THR:HG23	1:A:413:ALA:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:587:ARG:O	1:A:590:PRO:HD2	2.02	0.60
1:A:517:MET:HE3	1:A:529:LEU:HD21	1.84	0.60
1:A:587:ARG:HG2	1:A:587:ARG:O	2.02	0.59
1:A:206:THR:HG23	1:A:323:ALA:HB2	1.84	0.59
1:A:481:ARG:HD3	1:B:485:MET:SD	2.43	0.59
1:A:363:ARG:NE	1:A:400:ILE:HD11	2.19	0.58
1:B:348:PRO:HD2	1:B:380:LYS:HE2	1.86	0.57
1:B:460:GLN:NE2	1:B:464:ARG:HH11	2.02	0.56
1:B:347:ILE:HD11	1:B:356:LEU:HD13	1.88	0.56
1:A:381:LEU:HD13	1:A:386:ILE:HB	1.88	0.55
1:B:402:THR:CG2	1:B:403:ASN:N	2.69	0.55
1:B:432:VAL:HG22	1:B:450:THR:HG22	1.89	0.55
1:A:381:LEU:HB2	1:A:386:ILE:HD13	1.88	0.54
1:A:239:GLY:HA3	1:A:252:ILE:HD11	1.90	0.54
1:A:363:ARG:NH2	1:A:421:ASP:OD2	2.42	0.53
1:A:209:GLY:HA2	4:A:866:HOH:O	2.09	0.52
1:A:552:GLN:H	1:A:552:GLN:CD	2.13	0.52
1:B:592:LEU:O	1:B:594:GLY:N	2.43	0.52
1:A:364:HIS:HD2	1:A:424:SER:OG	1.93	0.51
1:A:426:ILE:HG12	1:A:474:ARG:HB2	1.92	0.51
1:B:413:ALA:O	1:B:414:LEU:HB3	2.10	0.51
1:B:213:LYS:HD3	1:B:213:LYS:H	1.74	0.51
1:A:598:LEU:HD22	1:A:609:VAL:HG11	1.92	0.51
1:B:418:PHE:CG	1:B:419:THR:N	2.79	0.51
1:B:601:ARG:HH21	1:B:605:VAL:HB	1.75	0.51
1:A:583:LYS:O	1:A:586:ILE:HG12	2.11	0.50
1:A:441:ASP:O	1:A:601:ARG:NH1	2.42	0.50
1:A:613:HIS:HE1	1:A:615:ILE:HD12	1.75	0.50
1:B:202:LEU:HD21	1:B:214:VAL:HG21	1.95	0.49
1:A:592:LEU:O	1:A:594:GLY:N	2.46	0.48
1:B:548:SER:O	1:B:552:GLN:HG3	2.14	0.48
1:A:400:ILE:HG21	1:A:419:THR:HG21	1.96	0.48
1:B:575:PRO:HB2	1:B:577:SER:O	2.14	0.47
1:A:411:THR:OG1	1:A:412:ASP:N	2.46	0.47
1:A:462:ARG:HG3	1:A:473:TYR:CG	2.50	0.47
1:A:613:HIS:CE1	1:A:615:ILE:HD12	2.50	0.46
1:A:236:LEU:HD23	1:A:252:ILE:HG21	1.96	0.46
1:A:379:ALA:HA	1:A:382:VAL:HG13	1.96	0.46
1:B:566:THR:O	1:B:570:ARG:HG3	2.16	0.46
1:A:464:ARG:HA	1:A:467:ARG:HH11	1.81	0.46
1:B:400:ILE:HA	1:B:401:PRO:HD2	1.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LYS:HB2	1:A:213:LYS:HE3	1.55	0.45
1:A:330:THR:HG23	1:A:480:GLU:OE1	2.15	0.45
1:B:485:MET:HG3	1:B:524:VAL:HG23	1.96	0.45
1:A:206:THR:HG22	1:A:210:LYS:NZ	2.31	0.45
1:A:386:ILE:O	1:A:387:ASN:HB3	2.17	0.45
1:B:204:ALA:O	1:B:323:ALA:HA	2.17	0.45
1:B:373:LYS:NZ	4:B:1112:HOH:O	2.50	0.45
1:A:450:THR:O	1:B:526:GLN:NE2	2.47	0.45
1:A:210:LYS:NZ	1:A:290:ASP:OD1	2.42	0.44
1:A:347:ILE:HD12	1:A:354:ILE:O	2.17	0.44
1:A:616:THR:HG22	1:A:620:MET:HE3	1.98	0.44
1:A:616:THR:HG22	1:A:620:MET:CE	2.47	0.44
1:B:542:ILE:HD11	1:B:562:ALA:HB3	1.99	0.44
1:B:212:THR:OG1	1:B:213:LYS:N	2.49	0.44
1:A:427:ASP:OD2	1:A:473:TYR:OH	2.35	0.44
1:B:431:CYS:O	1:B:450:THR:HA	2.18	0.44
1:B:591:THR:O	1:B:593:HIS:N	2.50	0.44
1:B:357:GLU:H	1:B:357:GLU:HG3	1.43	0.43
1:B:535:VAL:O	1:B:539:LEU:HG	2.18	0.43
1:A:519:THR:HA	1:A:520:PRO:HD3	1.85	0.43
1:B:616:THR:O	1:B:620:MET:HG3	2.19	0.43
1:A:458:ARG:NH2	1:A:477:ALA:O	2.51	0.43
1:B:518:ASN:ND2	4:B:1110:HOH:O	2.40	0.42
1:A:390:ALA:HA	1:A:409:VAL:O	2.19	0.42
1:B:563:TYR:HE2	1:B:615:ILE:HD13	1.83	0.42
1:A:366:ILE:O	1:A:409:VAL:HA	2.20	0.42
1:A:574:PRO:HG3	1:A:592:LEU:HA	2.01	0.42
1:B:469:LYS:HG2	1:B:470:PRO:N	2.34	0.42
1:B:613:HIS:CE1	1:B:615:ILE:HD12	2.54	0.42
1:B:616:THR:HG22	1:B:620:MET:CE	2.49	0.42
1:B:380:LYS:HE3	1:B:380:LYS:HB3	1.77	0.42
1:B:218:TYR:O	1:B:223:TYR:HB2	2.19	0.42
1:B:305:THR:OG1	1:B:512:ARG:HD3	2.20	0.42
1:A:469:LYS:HE3	1:A:469:LYS:HB3	1.67	0.42
1:B:415:MET:HG2	1:B:464:ARG:HH12	1.85	0.41
1:B:480:GLU:H	1:B:480:GLU:HG2	1.68	0.41
1:B:190:PRO:HA	1:B:191:PRO:HD3	1.83	0.41
1:A:187:ASN:HB3	1:A:189:SER:O	2.21	0.41
1:B:415:MET:HG2	1:B:464:ARG:NH1	2.36	0.41
1:A:369:HIS:CD2	1:A:370:SER:HB3	2.56	0.41
1:A:516:TYR:OH	1:A:523:PRO:HD2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:VAL:HG22	1:B:265:ILE:HG12	2.03	0.40
1:B:586:ILE:O	1:B:588:LEU:N	2.54	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:ARG:NH1	1:B:195:GLN:O[3_545]	2.19	0.01

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	438/464 (94%)	412 (94%)	23 (5%)	3 (1%)	22	21
1	B	413/464 (89%)	388 (94%)	21 (5%)	4 (1%)	15	13
All	All	851/928 (92%)	800 (94%)	44 (5%)	7 (1%)	19	17

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	587	ARG
1	B	213	LYS
1	B	401	PRO
1	B	419	THR
1	B	593	HIS
1	A	593	HIS
1	A	400	ILE

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	366/382 (96%)	361 (99%)	5 (1%)	67	76
1	B	346/382 (91%)	338 (98%)	8 (2%)	50	59
All	All	712/764 (93%)	699 (98%)	13 (2%)	59	68

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	SER
1	A	382	VAL
1	A	397	VAL
1	A	587	ARG
1	A	600	TYR
1	B	212	THR
1	B	357	GLU
1	B	395	LEU
1	B	402	THR
1	B	481	ARG
1	B	588	LEU
1	B	591	THR
1	B	600	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	364	HIS
1	A	526	GLN
1	B	198	GLN
1	B	460	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
2	2SX	A	701	-	11,15,15	0.83	0	12,21,21	1.32	1 (8%)

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	2SX	A	701	-	-	1/2/4/4	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	2SX	BR1-C1-C2	2.32	122.68	119.30

There are no chirality outliers.

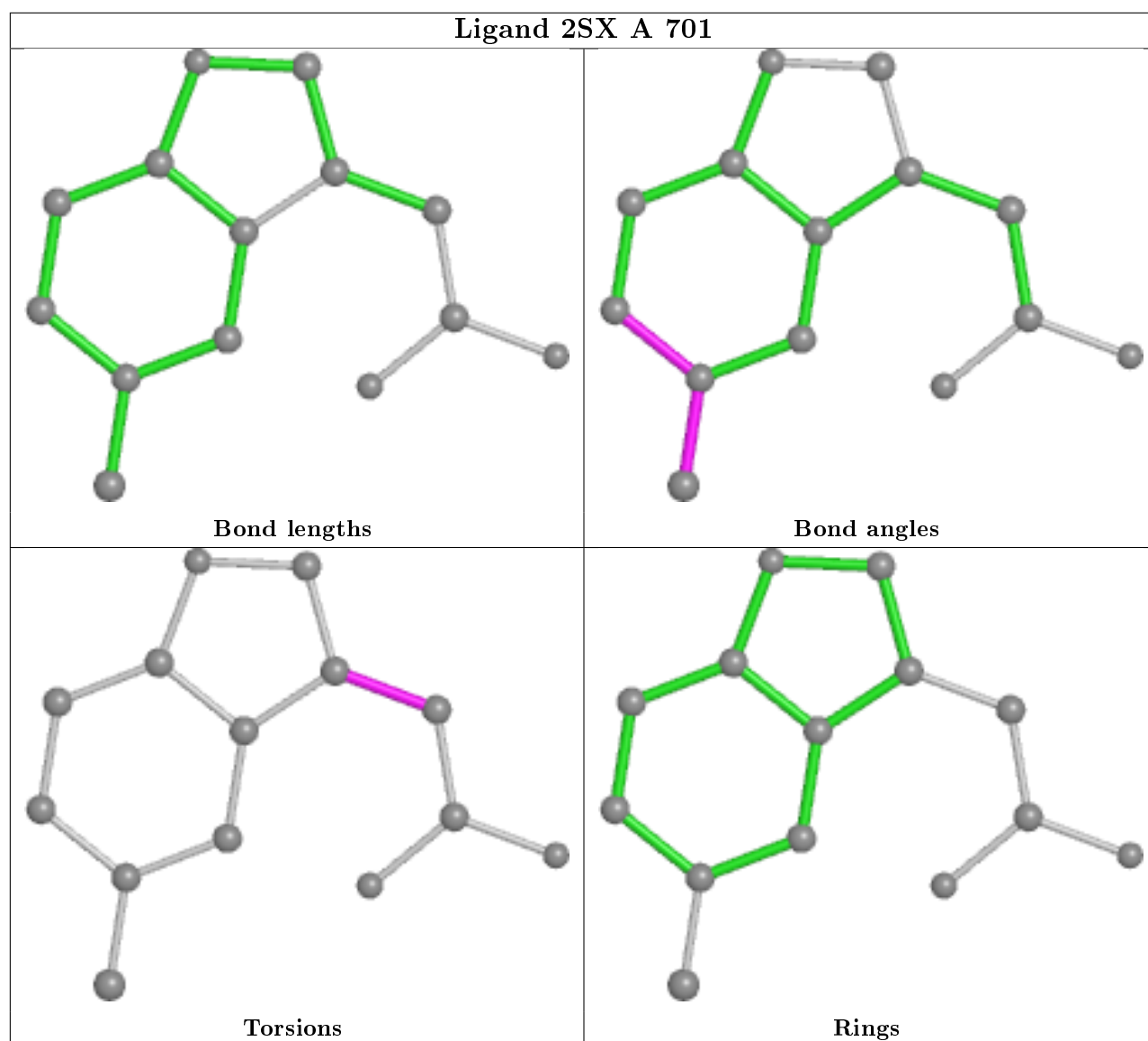
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	2SX	C12-C11-C9-C5

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	440/464 (94%)	0.78	48 (10%) 5 5	41, 59, 98, 114	0
1	B	420/464 (90%)	0.57	32 (7%) 13 15	39, 58, 94, 110	0
All	All	860/928 (92%)	0.68	80 (9%) 8 9	39, 59, 97, 114	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	420	GLY	8.4
1	A	418	PHE	7.9
1	A	383	ALA	7.7
1	A	382	VAL	7.2
1	B	401	PRO	7.1
1	A	592	LEU	6.8
1	B	592	LEU	6.2
1	B	263	SER	5.8
1	A	385	GLY	5.7
1	B	281	GLY	5.5
1	A	391	TYR	5.3
1	A	392	TYR	4.8
1	A	416	THR	4.7
1	B	212	THR	4.7
1	A	363	ARG	4.6
1	A	356	LEU	4.6
1	A	379	ALA	4.4
1	B	206	THR	4.4
1	A	281	GLY	4.3
1	B	220	ALA	4.3
1	A	423	ASP	4.2
1	A	206	THR	4.1
1	A	354	ILE	4.1
1	B	415	MET	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	358	VAL	3.9
1	B	396	ASP	3.9
1	B	236	LEU	3.9
1	B	188	SER	3.9
1	B	466	GLY	3.7
1	B	403	ASN	3.7
1	A	545	HIS	3.7
1	A	593	HIS	3.6
1	A	485[A]	MET	3.5
1	A	399	VAL	3.5
1	A	406	VAL	3.4
1	A	282	GLY	3.4
1	B	402	THR	3.3
1	B	246	HIS	3.3
1	A	588	LEU	3.3
1	A	591	THR	3.2
1	A	387	ASN	3.2
1	B	417	GLY	3.2
1	A	362	GLY	3.2
1	A	421	ASP	3.2
1	A	405	ASP	3.1
1	B	189	SER	3.1
1	A	630	VAL	2.9
1	A	377	LEU	2.9
1	B	264	PRO	2.9
1	A	361	GLY	2.9
1	B	586	ILE	2.8
1	B	591	THR	2.8
1	A	386	ILE	2.8
1	A	395	LEU	2.8
1	A	400	ILE	2.7
1	B	588	LEU	2.6
1	A	337	GLU	2.6
1	B	589	LYS	2.6
1	A	468	GLY	2.5
1	A	424	SER	2.5
1	A	380	LYS	2.4
1	B	392	TYR	2.4
1	B	579	ASP	2.4
1	B	400	ILE	2.4
1	B	219	ALA	2.4
1	A	396	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	384	LEU	2.4
1	A	381	LEU	2.4
1	B	242	MET	2.3
1	A	359	ILE	2.3
1	B	578	TRP	2.3
1	A	390	ALA	2.2
1	B	393	ARG	2.2
1	A	360	LYS	2.2
1	B	555	GLU	2.2
1	A	347	ILE	2.2
1	B	418	PHE	2.2
1	B	580	GLN	2.1
1	A	344	THR	2.0
1	A	333	HIS	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 carbohydrates in this entry.

6.4 Ligands [i](#)

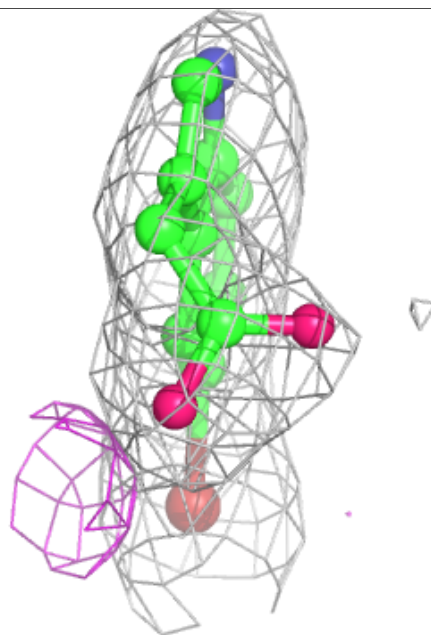
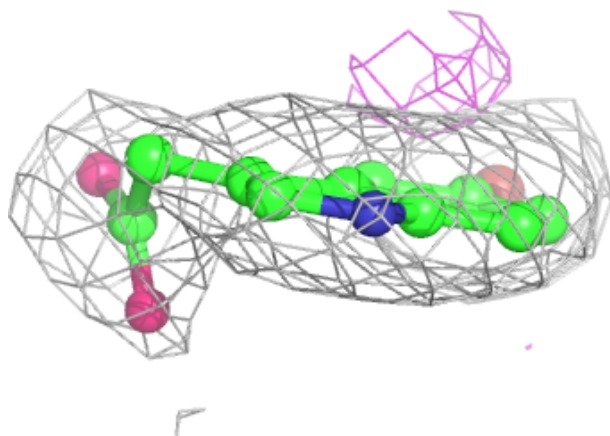
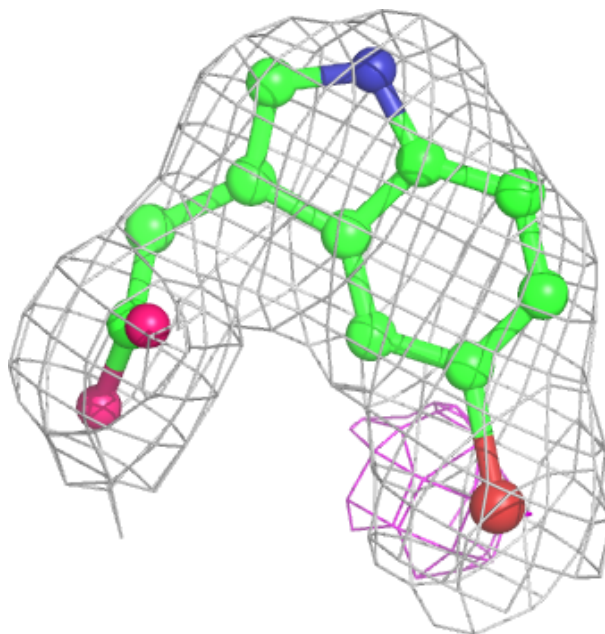
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CA	B	1001	1/1	0.62	0.12	76,76,76,76	0
2	2SX	A	701	14/14	0.92	0.15	55,61,68,115	0
3	CA	B	1002	1/1	0.93	0.09	81,81,81,81	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 2SX A 701:

$2mF_o-DF_c$ (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.