



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:01 AM UTC

PDB ID : 4JU2 / pdb_00004ju2
Title : Crystal structure of hcv ns5b polymerase in complex with compound 12
Authors : Coulombe, R.
Deposited on : 2013-03-24
Resolution : 2.70 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

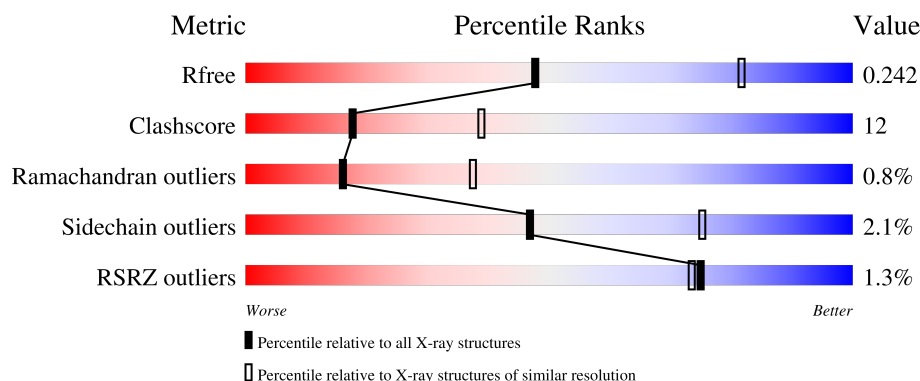
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	576	% 68% 27% ..
1	B	576	2% 69% 27% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9091 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	560	Total	C	N	O	S	0	0	0
			4362	2747	771	812	32			
1	B	560	Total	C	N	O	S	0	0	0
			4362	2747	771	812	32			

There are 12 discrepancies between the modelled and reference sequences:

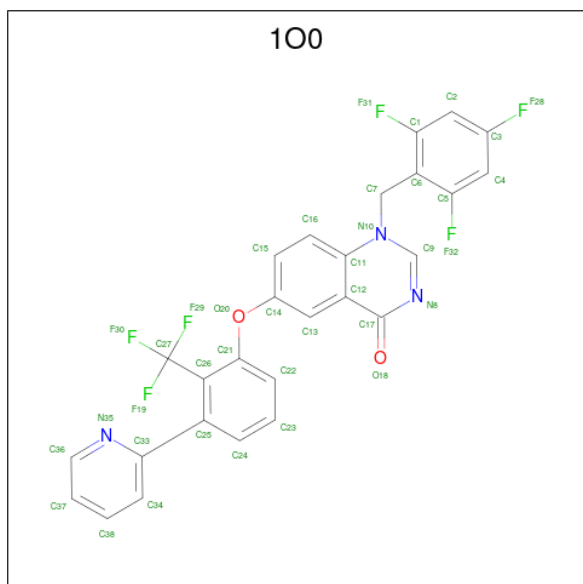
Chain	Residue	Modelled	Actual	Comment	Reference
A	571	HIS	-	expression tag	UNP O92972
A	572	HIS	-	expression tag	UNP O92972
A	573	HIS	-	expression tag	UNP O92972
A	574	HIS	-	expression tag	UNP O92972
A	575	HIS	-	expression tag	UNP O92972
A	576	HIS	-	expression tag	UNP O92972
B	571	HIS	-	expression tag	UNP O92972
B	572	HIS	-	expression tag	UNP O92972
B	573	HIS	-	expression tag	UNP O92972
B	574	HIS	-	expression tag	UNP O92972
B	575	HIS	-	expression tag	UNP O92972
B	576	HIS	-	expression tag	UNP O92972

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is 6-[3-(pyridin-2-yl)-2-(trifluoromethyl)phenoxy]-1-(2,4,6-trifluorobenzyl)quinazolin-4(1H)-one (CCD ID: 100) (formula: $C_{27}H_{15}F_6N_3O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			38	27	6	3	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	F	N	O	0	0
			38	27	6	3	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		
4	B	2	Total	Mg	0	0
			2	2		

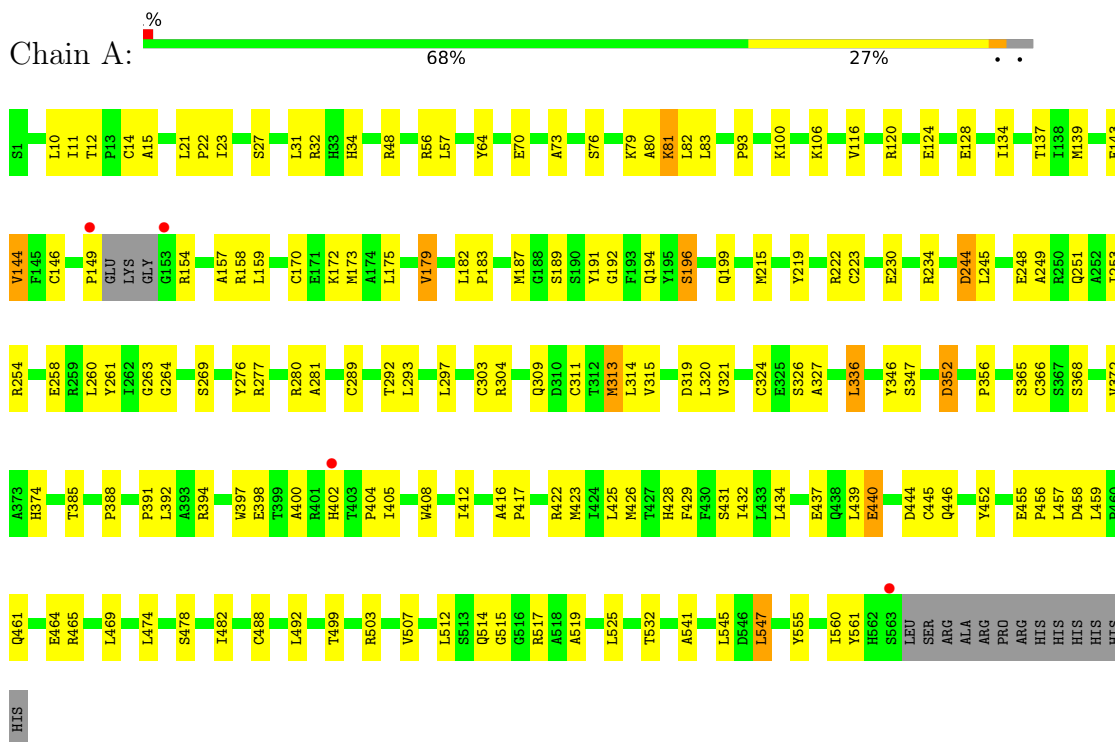
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	141	Total	O	0	0
			141	141		
5	B	134	Total	O	0	0
			134	134		

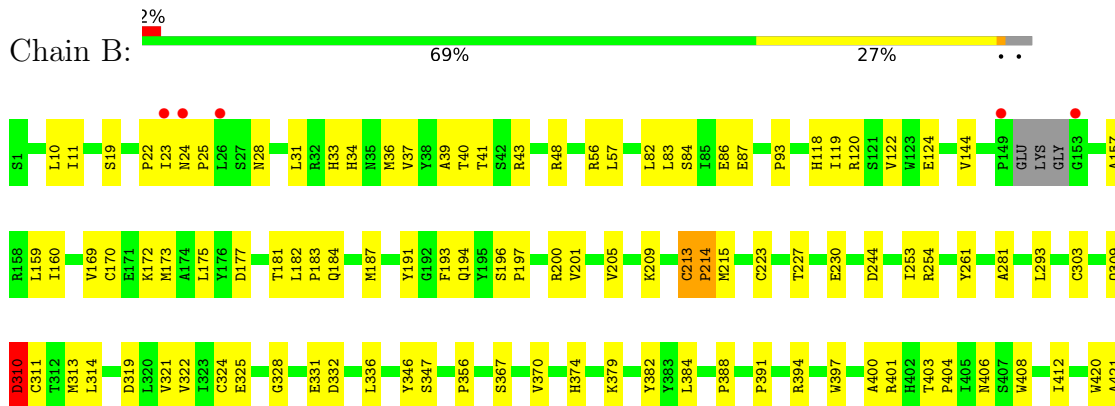
3 Residue-property plots [i](#)

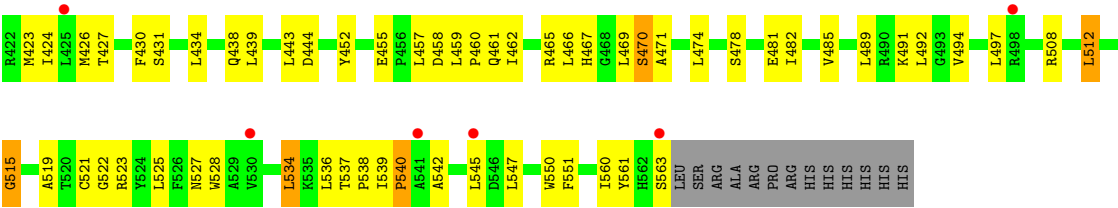
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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



• 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	105.54Å 107.74Å 135.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.70) 99.9 (50.00-2.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.69Å)	Xtriage
Refinement program	CNX 2002	Depositor
R, R_{free}	0.201 , 0.250 0.198 , 0.242	Depositor DCC
R_{free} test set	2170 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	51.8	Xtriage
Anisotropy	0.578	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.024 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9091	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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: MG, 100, GOL

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	0.44	0/4457	0.95	12/6049 (0.2%)
1	B	0.42	0/4457	0.96	16/6049 (0.3%)
All	All	0.43	0/8914	0.95	28/12098 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	SER	N-CA-C	-9.44	98.98	110.31
1	B	196	SER	N-CA-C	-7.87	100.87	110.31
1	B	537	THR	CA-C-N	7.27	128.93	119.84
1	B	537	THR	C-N-CA	7.27	128.93	119.84
1	A	100	LYS	N-CA-C	-7.23	104.40	112.57
1	B	261	TYR	N-CA-C	7.04	118.95	111.28
1	B	281	ALA	N-CA-C	-6.72	99.64	110.32
1	B	244	ASP	N-CA-C	-6.71	98.98	109.25
1	A	315	VAL	N-CA-C	6.66	117.88	108.36
1	A	425	LEU	N-CA-C	6.41	117.93	111.07
1	A	244	ASP	N-CA-C	-6.18	99.79	109.25
1	A	261	TYR	N-CA-C	6.03	117.93	111.36
1	A	352	ASP	N-CA-C	-5.79	101.58	110.32
1	B	213	CYS	CA-C-N	5.73	125.67	119.76
1	B	213	CYS	C-N-CA	5.73	125.67	119.76
1	B	430	PHE	N-CA-C	-5.70	105.59	112.54
1	A	192	GLY	N-CA-C	5.51	120.60	113.27
1	B	512	LEU	N-CA-C	-5.31	105.60	111.71
1	B	37	VAL	N-CA-C	5.30	116.93	108.87
1	B	420	TRP	N-CA-C	5.26	117.01	111.28
1	B	214	PRO	N-CA-C	5.23	119.29	111.14
1	B	200	ARG	N-CA-C	-5.22	105.48	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	A	429	PHE	N-CA-C	5.16	117.80	111.82
1	B	230	GLU	N-CA-C	-5.15	105.75	111.36
1	A	327	ALA	N-CA-C	-5.10	105.78	112.41
1	B	10	LEU	N-CA-C	5.08	117.41	110.55
1	A	525	LEU	N-CA-C	5.06	116.48	111.07
1	A	144	VAL	N-CA-C	5.01	115.80	108.58

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	4362	0	4374	104	0
1	B	4362	0	4374	114	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
3	A	38	0	15	1	0
3	B	38	0	15	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	141	0	0	8	0
5	B	134	0	0	4	0
All	All	9091	0	8794	217	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 12.

All (217) 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:14:CYS:HB2	1:A:139:MET:HE1	1.41	0.98
1:B:187:MET:HE2	1:B:293:LEU:HD23	1.62	0.81
1:A:248:GLU:HG3	5:A:3114:HOH:O	1.79	0.81
1:A:303:CYS:HG	1:A:311:CYS:HG	1.27	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:CYS:HG	1:B:311:CYS:HG	1.30	0.79
1:B:119:ILE:HD13	1:B:169:VAL:HG11	1.64	0.78
1:B:515:GLY:HA2	1:B:519:ALA:HB2	1.64	0.77
1:A:175:LEU:O	1:A:179:VAL:HG13	1.85	0.76
1:B:201:VAL:HG13	1:B:370:VAL:HG13	1.68	0.76
1:A:439:LEU:HB3	1:A:457:LEU:HD21	1.70	0.73
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.72	0.72
1:A:314:LEU:HB3	1:A:321:VAL:HG13	1.71	0.72
1:B:426:MET:HE1	1:B:525:LEU:HD22	1.72	0.71
1:A:22:PRO:HG2	1:A:400:ALA:HB1	1.74	0.70
1:A:503:ARG:O	1:A:507:VAL:HG23	1.92	0.70
1:B:439:LEU:HD23	1:B:457:LEU:HD21	1.73	0.70
1:A:394:ARG:O	1:A:398:GLU:HG3	1.91	0.69
1:B:33:HIS:HB2	1:B:492:LEU:O	1.92	0.69
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.23	0.68
1:A:440:GLU:HA	5:A:3143:HOH:O	1.92	0.68
1:A:499:THR:O	1:A:503:ARG:HG3	1.95	0.67
1:A:12:THR:HG1	1:A:269:SER:HG	1.43	0.67
1:B:336:LEU:HD23	1:B:356:PRO:HD3	1.77	0.67
1:A:154:ARG:HG3	1:A:154:ARG:HH11	1.60	0.67
1:B:408:TRP:O	1:B:412:ILE:HG13	1.95	0.66
1:B:314:LEU:HB3	1:B:321:VAL:HG13	1.78	0.66
1:B:382:TYR:HB2	5:B:780:HOH:O	1.95	0.66
1:B:93:PRO:HG3	1:B:561:TYR:HB2	1.78	0.65
1:B:515:GLY:CA	1:B:519:ALA:HB2	2.25	0.65
1:B:33:HIS:HB3	1:B:36:MET:HG2	1.77	0.65
1:B:434:LEU:HD23	1:B:439:LEU:HD11	1.80	0.64
1:B:367:SER:HB2	5:B:707:HOH:O	1.98	0.64
1:A:405:ILE:HG12	1:A:446:GLN:HE21	1.62	0.64
1:B:313:MET:SD	1:B:322:VAL:HG22	2.38	0.62
1:B:508:ARG:HG3	1:B:522:GLY:O	1.99	0.62
1:B:175:LEU:HD21	1:B:253:ILE:HG12	1.82	0.62
1:B:24:ASN:N	1:B:25:PRO:HD3	2.16	0.61
1:A:455:GLU:HB3	5:A:3143:HOH:O	1.99	0.61
1:A:223:CYS:HB3	5:A:3220:HOH:O	2.01	0.60
1:A:336:LEU:HD12	1:A:356:PRO:HD3	1.82	0.60
1:A:116:VAL:O	1:A:120:ARG:HG3	2.03	0.59
1:B:309:GLN:O	1:B:324:CYS:HB2	2.03	0.58
1:A:21:LEU:HD12	1:A:22:PRO:HD2	1.84	0.58
1:B:22:PRO:O	1:B:25:PRO:HG3	2.04	0.58
1:A:48:ARG:HG2	1:A:159:LEU:HG	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:422:ARG:HA	1:A:426:MET:SD	2.44	0.57
1:A:187:MET:HE2	1:A:293:LEU:HD23	1.86	0.57
1:A:461:GLN:HG2	1:A:541:ALA:HB3	1.86	0.57
1:A:263:GLY:HA2	1:A:277:ARG:NH1	2.20	0.57
1:A:515:GLY:HA2	1:A:519:ALA:HB2	1.87	0.56
1:B:144:VAL:HB	1:B:394:ARG:HG2	1.86	0.56
1:A:388:PRO:HG2	1:A:488:CYS:SG	2.45	0.56
1:A:172:LYS:HE3	1:A:560:ILE:HD13	1.87	0.56
1:B:465:ARG:HD3	1:B:542:ALA:O	2.06	0.56
1:B:328:GLY:HA3	1:B:331:GLU:OE1	2.05	0.56
1:A:93:PRO:HG3	1:A:561:TYR:HB2	1.87	0.55
1:B:444:ASP:HA	1:B:452:TYR:O	2.06	0.55
1:B:215:MET:HE2	1:B:332:ASP:HB3	1.89	0.55
1:B:421:ALA:O	1:B:426:MET:HG3	2.06	0.54
1:B:534:LEU:H	1:B:534:LEU:HD12	1.72	0.54
1:A:11:ILE:HD13	1:A:159:LEU:HD22	1.89	0.54
1:B:314:LEU:HB3	1:B:321:VAL:CG1	2.36	0.54
1:A:405:ILE:HG23	1:A:445:CYS:HA	1.90	0.54
1:B:187:MET:HE2	1:B:293:LEU:CD2	2.35	0.54
1:B:439:LEU:HB3	1:B:457:LEU:HD21	1.89	0.54
1:A:309:GLN:O	1:A:324:CYS:HB2	2.07	0.54
1:B:460:PRO:HB2	1:B:461:GLN:HE21	1.73	0.54
1:B:489:LEU:HD22	1:B:494:VAL:HG11	1.89	0.53
1:B:331:GLU:CD	1:B:331:GLU:H	2.16	0.53
1:A:532:THR:HG22	1:A:532:THR:O	2.09	0.53
1:A:21:LEU:HD23	1:A:34:HIS:HB2	1.91	0.53
1:A:80:ALA:HB3	1:A:245:LEU:HD23	1.90	0.53
1:A:478:SER:HB2	5:A:3188:HOH:O	2.09	0.52
1:B:48:ARG:HG2	1:B:159:LEU:HG	1.92	0.52
1:B:303:CYS:SG	1:B:313:MET:HE1	2.49	0.52
1:A:170:CYS:HA	1:A:173:MET:HE3	1.91	0.52
1:B:82:LEU:HD12	1:B:173:MET:O	2.09	0.52
1:B:120:ARG:NH1	1:B:124:GLU:HG3	2.25	0.52
1:A:79:LYS:HG3	1:A:244:ASP:HB3	1.91	0.52
1:B:461:GLN:H	1:B:461:GLN:NE2	2.09	0.51
1:B:36:MET:HE2	1:B:491:LYS:O	2.11	0.51
1:B:201:VAL:HG13	1:B:370:VAL:CG1	2.40	0.51
1:A:368:SER:HB2	1:A:385:THR:O	2.10	0.51
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.91	0.51
1:B:346:TYR:O	1:B:347:SER:HB3	2.11	0.51
1:A:437:GLU:HA	5:A:3234:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:LEU:HD13	5:B:703:HOH:O	2.11	0.50
1:A:146:CYS:SG	1:A:492:LEU:HD21	2.51	0.50
1:B:481:GLU:O	1:B:485:VAL:HG23	2.12	0.50
1:B:84:SER:OG	1:B:87:GLU:HG3	2.12	0.50
1:B:40:THR:HB	1:B:157:ALA:HB2	1.94	0.50
1:A:465:ARG:HG2	1:A:547:LEU:HD22	1.94	0.49
1:B:461:GLN:HG3	1:B:539:ILE:HG21	1.94	0.49
1:B:561:TYR:CE2	1:B:563:SER:HB3	2.47	0.49
1:B:427:THR:O	1:B:431:SER:HB2	2.12	0.49
1:B:397:TRP:CZ2	1:B:401:ARG:HD2	2.47	0.49
1:B:497:LEU:HD12	1:B:497:LEU:O	2.13	0.49
1:B:534:LEU:HD12	1:B:534:LEU:N	2.27	0.49
1:A:183:PRO:HG3	1:A:289:CYS:SG	2.52	0.49
1:A:388:PRO:C	1:A:391:PRO:HD2	2.37	0.49
1:A:465:ARG:NH1	1:A:545:LEU:O	2.46	0.49
1:A:515:GLY:CA	1:A:519:ALA:HB2	2.43	0.49
1:A:10:LEU:O	1:A:12:THR:HG23	2.13	0.49
1:A:144:VAL:HB	1:A:394:ARG:HG2	1.95	0.49
1:A:144:VAL:HG21	1:A:397:TRP:CG	2.48	0.49
1:A:249:ALA:O	1:A:253:ILE:HG13	2.13	0.48
1:B:388:PRO:C	1:B:391:PRO:HD2	2.38	0.48
1:B:31:LEU:HD12	1:B:31:LEU:O	2.12	0.48
1:B:56:ARG:O	1:B:57:LEU:HD23	2.14	0.48
1:B:170:CYS:HA	1:B:173:MET:HE3	1.96	0.48
1:B:182:LEU:HB3	1:B:183:PRO:HD3	1.96	0.48
1:A:191:TYR:O	1:A:194:GLN:HG2	2.13	0.47
1:B:215:MET:SD	1:B:215:MET:C	2.97	0.47
1:B:527:ASN:ND2	1:B:536:LEU:HD11	2.28	0.47
1:A:313:MET:HG3	1:A:320:LEU:HD11	1.95	0.47
1:B:406:ASN:ND2	1:B:443:LEU:HB3	2.29	0.47
1:B:523:ARG:O	1:B:527:ASN:HB2	2.15	0.47
1:A:405:ILE:HG12	1:A:446:GLN:NE2	2.28	0.47
1:B:197:PRO:O	1:B:201:VAL:HG23	2.15	0.47
1:A:48:ARG:HD3	1:A:157:ALA:O	2.14	0.46
1:B:460:PRO:HD2	1:B:461:GLN:HE22	1.80	0.46
1:A:405:ILE:N	1:A:405:ILE:HD12	2.30	0.46
1:B:28:ASN:HA	1:B:34:HIS:NE2	2.29	0.46
1:B:11:ILE:HD13	1:B:159:LEU:HD22	1.98	0.46
1:B:374:HIS:HA	1:B:379:LYS:O	2.16	0.46
1:A:478:SER:O	1:A:482:ILE:HG13	2.16	0.46
1:A:196:SER:OG	1:A:199:GLN:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:TYR:CZ	1:B:193:PHE:HB2	2.50	0.46
1:A:134:ILE:HD12	1:A:260:LEU:HB2	1.97	0.45
1:A:254:ARG:NH2	1:A:258:GLU:HG2	2.31	0.45
1:A:440:GLU:H	1:A:440:GLU:CD	2.23	0.45
1:B:191:TYR:CE2	1:B:193:PHE:HB2	2.51	0.45
1:A:428:HIS:O	1:A:432:ILE:HG13	2.16	0.45
1:B:309:GLN:HB3	1:B:325:GLU:HB3	1.98	0.45
1:A:81:LYS:HE2	1:A:83:LEU:CD1	2.46	0.45
1:A:189:SER:HA	1:A:194:GLN:HE22	1.82	0.45
1:A:512:LEU:C	1:A:514:GLN:H	2.25	0.45
1:B:201:VAL:CG2	1:B:384:LEU:HB2	2.47	0.45
1:B:331:GLU:CD	1:B:331:GLU:N	2.75	0.45
1:B:374:HIS:O	1:B:474:LEU:HA	2.16	0.45
1:A:455:GLU:HB2	1:A:458:ASP:OD2	2.17	0.45
1:A:81:LYS:HE2	1:A:83:LEU:HD12	2.00	0.44
1:A:321:VAL:HB	1:A:365:SER:HB3	1.99	0.44
1:A:222:ARG:HG2	1:A:222:ARG:HH11	1.82	0.44
1:A:251:GLN:NE2	1:B:254:ARG:HG2	2.33	0.44
1:B:160:ILE:O	1:B:160:ILE:HG13	2.17	0.44
1:B:426:MET:CE	1:B:525:LEU:HD22	2.44	0.44
1:A:154:ARG:HG3	1:A:154:ARG:NH1	2.29	0.44
1:A:541:ALA:O	1:A:545:LEU:HG	2.18	0.43
1:B:466:LEU:HD22	1:B:551:PHE:HE2	1.83	0.43
1:B:469:LEU:C	1:B:471:ALA:H	2.27	0.43
1:B:508:ARG:NH1	1:B:512:LEU:HD11	2.33	0.43
1:B:172:LYS:HE3	1:B:560:ILE:HD13	2.00	0.43
1:A:57:LEU:HD23	1:A:57:LEU:C	2.43	0.43
1:A:374:HIS:O	1:A:474:LEU:HA	2.19	0.43
1:A:423:MET:HE1	3:A:3002:1O0:H12	2.01	0.43
1:B:455:GLU:HB2	1:B:458:ASP:OD2	2.18	0.43
1:A:280:ARG:HG3	1:A:281:ALA:O	2.18	0.43
1:B:177:ASP:O	1:B:181:THR:HG23	2.18	0.43
1:B:536:LEU:N	1:B:536:LEU:HD12	2.33	0.43
1:B:118:HIS:O	1:B:122:VAL:HG23	2.19	0.43
1:B:423:MET:HE3	1:B:424:ILE:HD11	2.00	0.43
1:A:456:PRO:O	1:A:459:LEU:HG	2.19	0.43
1:A:555:TYR:CD1	1:A:560:ILE:HG13	2.54	0.43
1:A:416:ALA:N	1:A:417:PRO:CD	2.82	0.42
1:A:434:LEU:CD2	1:A:439:LEU:HD11	2.48	0.42
1:B:478:SER:O	1:B:482:ILE:HG13	2.19	0.42
1:A:457:LEU:HD12	1:A:517:ARG:NH2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:GLU:OE2	1:A:158:ARG:NH2	2.53	0.42
1:B:184:GLN:HB2	5:B:803:HOH:O	2.19	0.42
1:A:222:ARG:HD3	5:A:3140:HOH:O	2.19	0.42
1:B:539:ILE:HG23	1:B:540:PRO:HD2	2.01	0.42
1:A:263:GLY:HA3	1:A:277:ARG:O	2.20	0.42
1:A:264:GLY:HA2	1:A:276:TYR:CZ	2.55	0.42
1:A:392:LEU:HD23	1:A:392:LEU:HA	1.87	0.42
1:B:23:ILE:O	1:B:24:ASN:HB2	2.20	0.42
1:B:33:HIS:HB3	1:B:36:MET:CG	2.46	0.42
1:B:86:GLU:OE1	1:B:86:GLU:N	2.52	0.42
1:A:444:ASP:HA	1:A:452:TYR:O	2.19	0.42
1:A:56:ARG:HD3	1:A:56:ARG:N	2.35	0.41
1:A:230:GLU:O	1:A:234:ARG:HG3	2.20	0.41
1:A:402:HIS:CD2	5:A:3157:HOH:O	2.72	0.41
1:B:184:GLN:HE22	1:B:194:GLN:CD	2.27	0.41
1:B:397:TRP:CE2	1:B:401:ARG:HD2	2.55	0.41
1:A:215:MET:HB2	1:A:326:SER:HB2	2.01	0.41
1:A:404:PRO:HB2	1:A:405:ILE:HD12	2.01	0.41
1:A:464:GLU:OE1	1:A:469:LEU:HD21	2.20	0.41
1:B:223:CYS:O	1:B:227:THR:HG23	2.20	0.41
1:B:403:THR:HA	1:B:404:PRO:HD3	1.94	0.41
1:B:547:LEU:O	1:B:550:TRP:HB2	2.20	0.41
1:A:73:ALA:O	1:A:76:SER:HB2	2.21	0.41
1:B:521:CYS:O	1:B:525:LEU:HB2	2.20	0.41
1:A:346:TYR:O	1:A:347:SER:HB3	2.20	0.41
1:B:11:ILE:HG22	1:B:11:ILE:O	2.20	0.41
1:A:31:LEU:HB2	1:A:492:LEU:O	2.21	0.41
1:A:124:GLU:O	1:A:128:GLU:HG3	2.20	0.41
1:A:187:MET:HE1	1:A:292:THR:HG22	2.02	0.41
1:B:19:SER:HA	1:B:39:ALA:HB3	2.01	0.41
1:B:197:PRO:HB2	1:B:467:HIS:HE1	1.86	0.41
1:B:470:SER:O	1:B:474:LEU:HG	2.19	0.41
1:A:32:ARG:N	1:A:32:ARG:HD2	2.36	0.41
1:A:64:TYR:CZ	1:A:297:LEU:HD21	2.56	0.41
1:A:431:SER:HB2	1:A:507:VAL:HG21	2.03	0.41
1:B:191:TYR:O	1:B:194:GLN:HG2	2.21	0.41
1:B:205:VAL:HG13	1:B:209:LYS:HE3	2.02	0.41
1:B:213:CYS:HA	1:B:214:PRO:HD2	1.92	0.41
1:B:459:LEU:N	1:B:460:PRO:CD	2.83	0.41
1:B:41:THR:C	1:B:43:ARG:H	2.28	0.41
1:B:458:ASP:O	1:B:462:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:469:LEU:C	1:B:471:ALA:N	2.77	0.41
1:A:408:TRP:O	1:A:412:ILE:HG13	2.22	0.40
1:A:70:GLU:CD	1:A:304:ARG:HH22	2.29	0.40
1:B:197:PRO:HB2	1:B:467:HIS:CE1	2.56	0.40
1:B:83:LEU:HB2	1:B:173:MET:HA	2.04	0.40
1:B:310:ASP:O	1:B:324:CYS:HA	2.22	0.40
1:B:314:LEU:HD12	1:B:314:LEU:HA	1.75	0.40
1:B:523:ARG:HH11	1:B:523:ARG:HG3	1.85	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	556/576 (96%)	526 (95%)	29 (5%)	1 (0%)	43	68
1	B	556/576 (96%)	513 (92%)	35 (6%)	8 (1%)	9	23
All	All	1112/1152 (96%)	1039 (93%)	64 (6%)	9 (1%)	16	37

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	538	PRO
1	B	540	PRO
1	A	15	ALA
1	B	400	ALA
1	B	528	TRP
1	B	438	GLN
1	B	470	SER
1	B	310	ASP
1	B	515	GLY

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	477/491 (97%)	461 (97%)	16 (3%)	32	62
1	B	477/491 (97%)	473 (99%)	4 (1%)	73	88
All	All	954/982 (97%)	934 (98%)	20 (2%)	47	75

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ILE
1	A	27	SER
1	A	81	LYS
1	A	106	LYS
1	A	137	THR
1	A	149	PRO
1	A	179	VAL
1	A	182	LEU
1	A	313	MET
1	A	319	ASP
1	A	336	LEU
1	A	352	ASP
1	A	366	CYS
1	A	372	VAL
1	A	440	GLU
1	A	547	LEU
1	B	310	ASP
1	B	319	ASP
1	B	534	LEU
1	B	545	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	194	GLN

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Mol	Chain	Res	Type
1	A	273	ASN
1	A	406	ASN
1	A	436	GLN
1	A	438	GLN
1	A	446	GLN
1	A	502	HIS
1	A	514	GLN
1	A	527	ASN
1	A	544	GLN
1	A	562	HIS
1	B	117	ASN
1	B	184	GLN
1	B	206	ASN
1	B	273	ASN
1	B	330	GLN
1	B	402	HIS
1	B	406	ASN
1	B	438	GLN
1	B	461	GLN
1	B	467	HIS

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 8 ligands modelled in this entry, 4 are 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	$\# Z > 2$
2	GOL	B	601	-	5,5,5	0.37	0	5,5,5	0.24	0
3	100	A	3002	-	42,42,42	2.17	16 (38%)	59,62,62	1.66	13 (22%)
2	GOL	A	3001	-	5,5,5	0.54	0	5,5,5	0.24	0
3	100	B	602	-	42,42,42	2.50	20 (47%)	59,62,62	1.73	16 (27%)

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	GOL	B	601	-	-	4/4/4/4	-
3	100	A	3002	-	-	4/18/18/18	0/5/5/5
2	GOL	A	3001	-	-	2/4/4/4	-
3	100	B	602	-	-	4/18/18/18	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	100	C25-C26	6.88	1.49	1.40
3	A	3002	100	C25-C26	5.79	1.47	1.40
3	B	602	100	C11-N10	4.88	1.46	1.40
3	B	602	100	C21-C26	4.31	1.47	1.40
3	A	3002	100	C11-N10	4.15	1.45	1.40
3	A	3002	100	C21-C26	4.03	1.46	1.40
3	B	602	100	C6-C5	3.93	1.44	1.38
3	B	602	100	C6-C1	3.54	1.43	1.38
3	B	602	100	C12-C11	3.46	1.46	1.41
3	A	3002	100	C12-C11	3.35	1.46	1.41
3	A	3002	100	C6-C5	3.22	1.43	1.38
3	B	602	100	C2-C3	3.09	1.42	1.37
3	B	602	100	C13-C12	2.96	1.44	1.39
3	B	602	100	C4-C3	2.90	1.42	1.37
3	B	602	100	C2-C1	2.90	1.42	1.37
3	B	602	100	C24-C25	2.88	1.44	1.40
3	A	3002	100	C24-C25	2.81	1.44	1.40
3	B	602	100	C15-C14	2.81	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	3002	100	C13-C12	2.76	1.44	1.39
3	B	602	100	C9-N10	2.72	1.40	1.36
3	A	3002	100	C9-N10	2.71	1.40	1.36
3	B	602	100	C7-N10	2.68	1.51	1.47
3	B	602	100	C4-C5	2.67	1.42	1.37
3	A	3002	100	C6-C1	2.66	1.42	1.38
3	A	3002	100	C2-C1	2.60	1.42	1.37
3	A	3002	100	C2-C3	2.53	1.41	1.37
3	A	3002	100	C15-C14	2.49	1.43	1.38
3	A	3002	100	C7-N10	2.43	1.50	1.47
3	B	602	100	C16-C11	2.42	1.43	1.39
3	B	602	100	C16-C15	2.38	1.42	1.38
3	B	602	100	C33-N35	2.19	1.39	1.35
3	A	3002	100	C4-C5	2.13	1.41	1.37
3	B	602	100	C36-N35	2.13	1.39	1.34
3	B	602	100	C22-C21	2.07	1.43	1.39
3	A	3002	100	C22-C21	2.06	1.43	1.39
3	A	3002	100	C16-C15	2.03	1.42	1.38

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	100	C16-C11-N10	6.68	125.91	120.97
3	A	3002	100	C16-C11-N10	6.02	125.42	120.97
3	A	3002	100	C2-C1-C6	-4.03	121.45	124.45
3	B	602	100	C4-C5-C6	-3.77	121.64	124.45
3	B	602	100	C2-C1-C6	-3.50	121.85	124.45
3	B	602	100	C13-C12-C11	3.25	122.59	118.76
3	A	3002	100	C4-C5-C6	-3.23	122.05	124.45
3	A	3002	100	C13-C12-C11	3.05	122.36	118.76
3	A	3002	100	C27-C26-C25	2.86	124.21	120.62
3	B	602	100	C11-C12-C17	-2.79	117.67	119.99
3	A	3002	100	C11-C12-C17	-2.78	117.68	119.99
3	B	602	100	C33-C25-C26	2.58	126.78	122.94
3	B	602	100	C5-C4-C3	2.54	119.37	116.67
3	A	3002	100	N10-C9-N8	2.53	128.46	126.34
3	B	602	100	C16-C11-C12	-2.52	115.30	119.16
3	B	602	100	N10-C9-N8	2.46	128.40	126.34
3	A	3002	100	C33-C25-C26	2.39	126.50	122.94
3	B	602	100	F32-C5-C6	2.34	120.56	117.64
3	B	602	100	C27-C26-C25	2.32	123.54	120.62
3	A	3002	100	C16-C11-C12	-2.32	115.60	119.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	100	C4-C3-C2	-2.24	120.77	123.50
3	B	602	100	C1-C2-C3	2.24	119.06	116.67
3	A	3002	100	C1-C2-C3	2.24	119.06	116.67
3	B	602	100	C24-C25-C33	-2.23	114.69	119.56
3	A	3002	100	C24-C25-C33	-2.14	114.87	119.56
3	A	3002	100	C5-C6-C1	2.10	117.37	114.54
3	A	3002	100	C5-C4-C3	2.08	118.88	116.67
3	B	602	100	C5-C6-C1	2.05	117.30	114.54
3	B	602	100	O18-C17-C12	-2.02	118.08	121.33

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	3001	GOL	O1-C1-C2-C3
2	B	601	GOL	O1-C1-C2-C3
2	B	601	GOL	C1-C2-C3-O3
2	B	601	GOL	O2-C2-C3-O3
3	B	602	100	C24-C25-C33-N35
3	A	3002	100	C24-C25-C33-C34
3	B	602	100	C24-C25-C33-C34
3	A	3002	100	C24-C25-C33-N35
2	B	601	GOL	O1-C1-C2-O2
3	A	3002	100	C26-C25-C33-C34
3	B	602	100	C26-C25-C33-C34
2	A	3001	GOL	O1-C1-C2-O2
3	A	3002	100	C26-C25-C33-N35
3	B	602	100	C26-C25-C33-N35

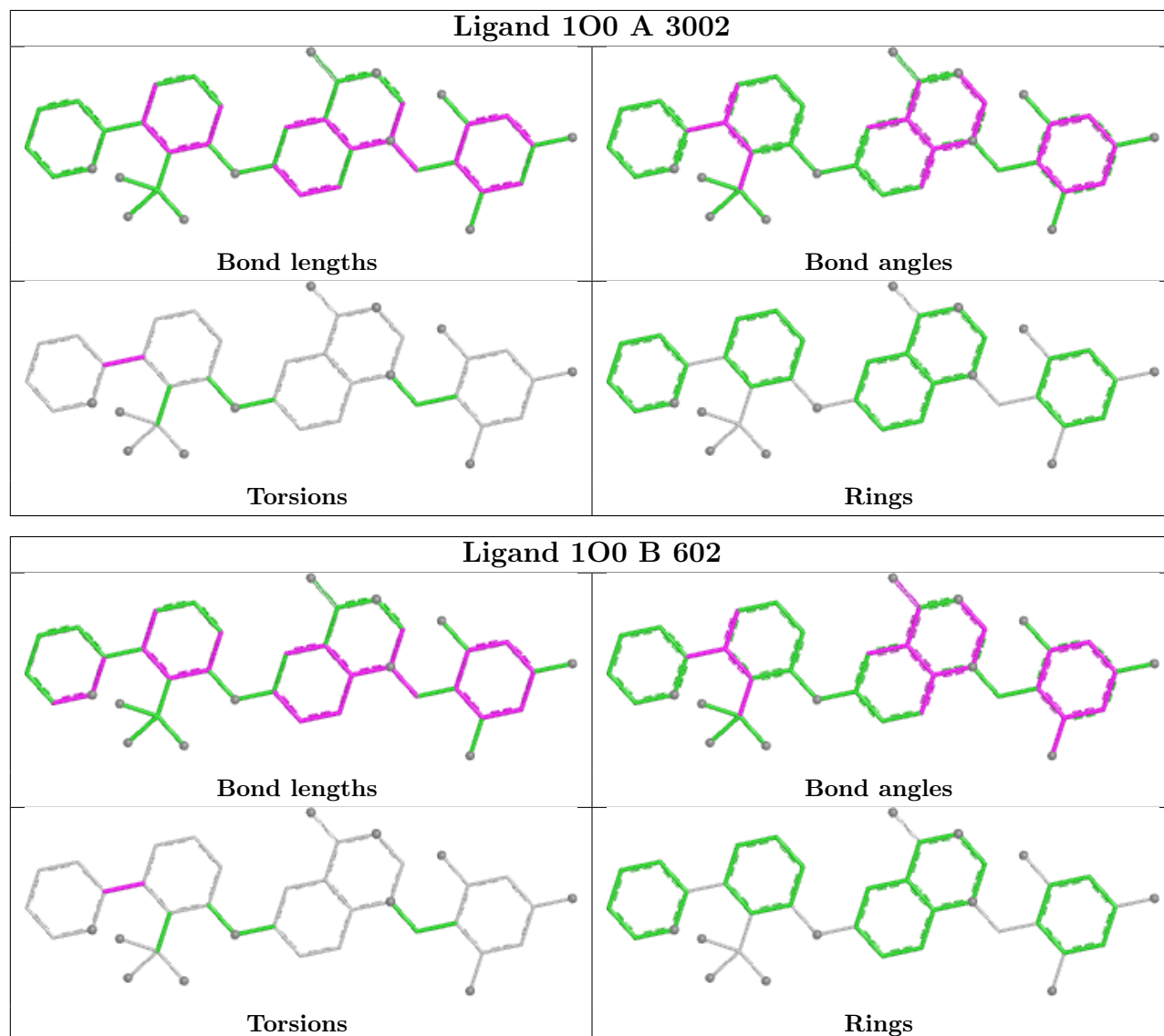
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	3002	100	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	560/576 (97%)	-0.30	4 (0%) 84 83	33, 50, 78, 98	0
1	B	560/576 (97%)	-0.04	11 (1%) 65 62	32, 63, 102, 117	0
All	All	1120/1152 (97%)	-0.17	15 (1%) 75 73	32, 55, 94, 117	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	149	PRO	4.5
1	B	541	ALA	4.0
1	B	23	ILE	3.7
1	A	149	PRO	3.7
1	B	26	LEU	2.9
1	B	498	ARG	2.8
1	B	563	SER	2.4
1	B	545	LEU	2.4
1	A	402	HIS	2.4
1	B	153	GLY	2.2
1	A	563	SER	2.2
1	B	425	LEU	2.2
1	A	153	GLY	2.1
1	B	530	VAL	2.1
1	B	24	ASN	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

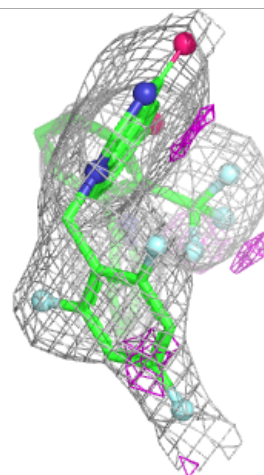
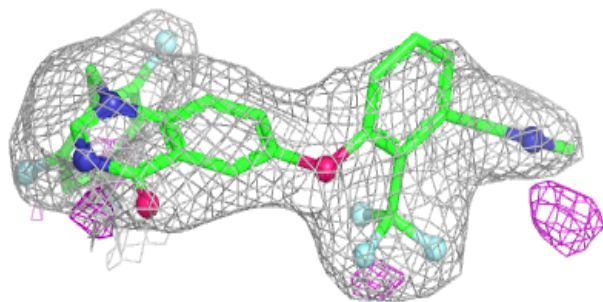
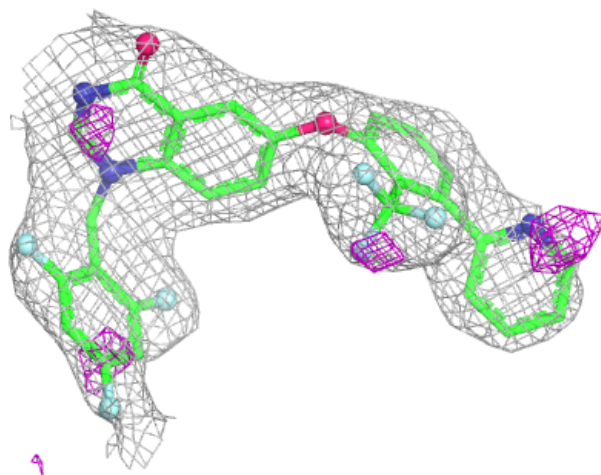
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	B	603	1/1	0.83	0.25	77,77,77,77	0
4	MG	B	604	1/1	0.84	0.18	83,83,83,83	0
4	MG	A	3004	1/1	0.87	0.15	66,66,66,66	0
3	1O0	B	602	38/38	0.90	0.11	62,68,78,78	0
4	MG	A	3003	1/1	0.91	0.12	65,65,65,65	0
2	GOL	B	601	6/6	0.92	0.17	80,80,81,81	0
3	1O0	A	3002	38/38	0.93	0.10	50,53,57,57	0
2	GOL	A	3001	6/6	0.93	0.12	60,65,67,68	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.

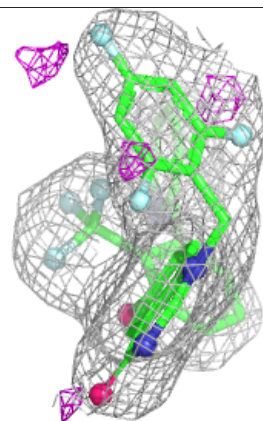
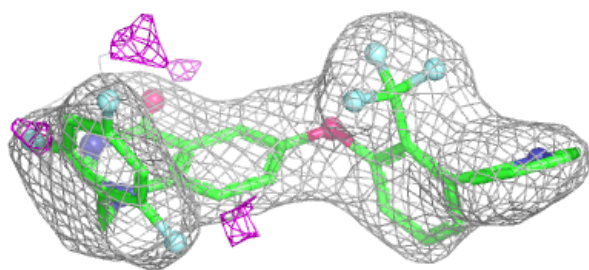
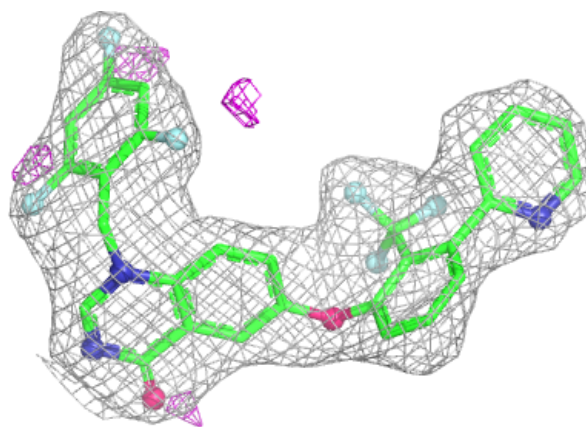
Electron density around 1O0 B 602:

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



Electron density around 1O0 A 3002:

$2mF_o-DF_c$ (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.