



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:43 PM UTC

PDB ID : 8B9P / pdb_00008b9p
Title : ACE2 in complex with bicyclic peptide inhibitor
Authors : Brear, P.; Lulla, A.; Harman, M.; Dods, R.; Chen, L.; Bezerra, G.; Demydchuk, Y.; Stanway, S.; Hyvonen, M.
Deposited on : 2022-10-06
Resolution : 2.11 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

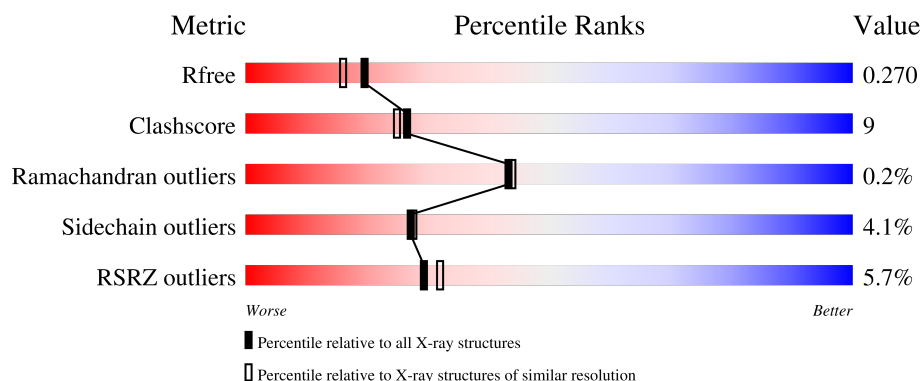
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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	609	4% 78% 19% .
1	B	609	7% 73% 24% .
2	C	18	11% 72% 22% 6%
2	D	18	6% 72% 22% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10085 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 Processed angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4862	3111	805	917	29			
1	B	595	Total	C	N	O	S	0	0	0
			4856	3108	804	915	29			

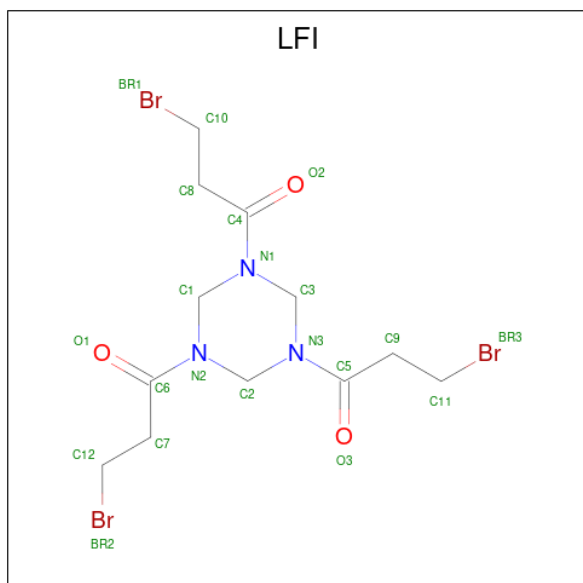
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	GLY	-	expression tag	UNP Q9BYF1
A	616	SER	-	expression tag	UNP Q9BYF1
A	617	SER	-	expression tag	UNP Q9BYF1
A	618	PRO	-	expression tag	UNP Q9BYF1
A	619	HIS	-	expression tag	UNP Q9BYF1
A	620	HIS	-	expression tag	UNP Q9BYF1
A	621	HIS	-	expression tag	UNP Q9BYF1
A	622	HIS	-	expression tag	UNP Q9BYF1
A	623	HIS	-	expression tag	UNP Q9BYF1
A	624	HIS	-	expression tag	UNP Q9BYF1
A	625	HIS	-	expression tag	UNP Q9BYF1
A	626	HIS	-	expression tag	UNP Q9BYF1
B	18	GLY	-	expression tag	UNP Q9BYF1
B	616	SER	-	expression tag	UNP Q9BYF1
B	617	SER	-	expression tag	UNP Q9BYF1
B	618	PRO	-	expression tag	UNP Q9BYF1
B	619	HIS	-	expression tag	UNP Q9BYF1
B	620	HIS	-	expression tag	UNP Q9BYF1
B	621	HIS	-	expression tag	UNP Q9BYF1
B	622	HIS	-	expression tag	UNP Q9BYF1
B	623	HIS	-	expression tag	UNP Q9BYF1
B	624	HIS	-	expression tag	UNP Q9BYF1
B	625	HIS	-	expression tag	UNP Q9BYF1
B	626	HIS	-	expression tag	UNP Q9BYF1

- Molecule 2 is a protein called ALA-CYS-GLY-ARG-GLN-PHE-CYS-HIS-THR-LEU-MET-PRO-ARG-HIS-LEU-CYS-ALA-NH₂.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	18	Total	C	N	O	S	0	0	1
			133	81	29	19	4			
2	D	18	Total	C	N	O	S	0	0	1
			133	81	29	19	4			

- Molecule 3 is 1-[3,5-bis(3-bromanylpropanoyl)-1,3,5-triazinan-1-yl]-3-bromanyl-propan-1-one (CCD ID: LFI) (formula: C₁₂H₁₈Br₃N₃O₃) (labeled as "Ligand of Interest" by depositor).



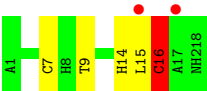
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			18	12	3	3		
3	D	1	Total	C	N	O	0	0
			18	12	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total	O	0	0
			41	41		
4	B	19	Total	O	0	0
			19	19		
4	C	4	Total	O	0	0
			4	4		
4	D	1	Total	O	0	0
			1	1		



• Molecule 2: ALA-CYS-GLY-ARG-GLN-PHE-CYS-HIS-THR-LEU-MET-PRO-ARG-HIS-LEU-CYS-ALA-NH2



• Molecule 2: ALA-CYS-GLY-ARG-GLN-PHE-CYS-HIS-THR-LEU-MET-PRO-ARG-HIS-LEU-CYS-ALA-NH2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.89Å 76.76Å 115.05Å 90.00° 99.93° 90.00°	Depositor
Resolution (Å)	66.69 – 2.11 66.69 – 2.11	Depositor EDS
% Data completeness (in resolution range)	97.2 (66.69-2.11) 97.2 (66.69-2.11)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.225 , 0.295 (Not available) , 0.270	Depositor DCC
R_{free} test set	3532 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	51.3	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10085	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 4.62% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NH2, LFI

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	1.00	0/4999	1.40	7/6792 (0.1%)
1	B	1.00	0/4993	1.38	5/6784 (0.1%)
2	C	1.46	1/135 (0.7%)	2.23	4/180 (2.2%)
2	D	1.09	0/135	2.07	3/180 (1.7%)
All	All	1.01	1/10262 (0.0%)	1.41	19/13936 (0.1%)

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
2	C	0	2
2	D	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	16	CYS	C-N	12.10	1.50	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	16	CYS	CA-C-N	16.47	151.35	121.70
2	C	16	CYS	C-N-CA	16.47	151.35	121.70
2	D	16	CYS	O-C-N	-14.88	102.26	122.19
2	D	16	CYS	CA-C-N	11.03	141.55	121.70
2	D	16	CYS	C-N-CA	11.03	141.55	121.70
1	A	269	ASP	CA-CB-CG	8.73	121.33	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	605	GLY	CA-C-O	-6.64	117.52	122.37
1	A	213	ASP	CA-CB-CG	6.34	118.94	112.60
1	A	352	GLY	CA-C-O	-6.18	118.00	122.45
1	B	117	ASN	CA-CB-CG	-6.01	106.59	112.60
2	C	15	LEU	N-CA-C	-5.72	106.01	112.87
1	A	221	GLN	N-CA-C	-5.57	105.30	111.71
1	A	350	ASP	CA-CB-CG	5.40	118.00	112.60
2	C	9	THR	CA-CB-OG1	-5.31	101.63	109.60
1	B	376	MET	N-CA-C	-5.17	105.65	111.28
1	B	437	ASN	CA-C-N	5.12	127.09	120.44
1	B	437	ASN	C-N-CA	5.12	127.09	120.44
1	B	549	GLU	N-CA-C	-5.06	105.45	110.97
1	A	394	ASN	CA-C-O	-5.00	115.38	120.89

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	16	CYS	Mainchain,Peptide
2	D	16	CYS	Peptide

5.2 Too-close contacts [i](#)

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	4862	0	4639	79	0
1	B	4856	0	4634	105	0
2	C	133	0	129	2	0
2	D	133	0	129	3	0
3	C	18	0	0	0	0
3	D	18	0	0	0	0
4	A	41	0	0	0	0
4	B	19	0	0	0	0
4	C	4	0	0	0	0
4	D	1	0	0	0	0
All	All	10085	0	9531	182	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 9.

All (182) 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:573:VAL:HG13	1:A:574:VAL:HG13	1.55	0.87
1:B:431:ASP:HB2	1:B:434:THR:HG23	1.63	0.80
1:B:311:ALA:HA	1:B:373:HIS:CE1	2.15	0.80
1:B:419:LYS:HZ3	1:B:428:PHE:HB3	1.48	0.79
1:B:538:PRO:HD2	1:B:541:LYS:HD3	1.67	0.76
1:A:345:HIS:O	1:A:347:THR:HG22	1.88	0.73
1:A:85:LEU:HB3	1:A:94:LYS:HE3	1.72	0.70
1:B:309:LYS:HD2	1:B:328:TRP:CH2	2.28	0.69
1:B:526:GLN:HA	1:B:526:GLN:OE1	1.91	0.69
1:A:245:ARG:HB2	1:A:262:LEU:HD21	1.74	0.68
1:A:309:LYS:HD2	1:A:328:TRP:CH2	2.29	0.67
1:A:168:TRP:CZ3	1:A:172:VAL:HG21	2.30	0.67
1:B:517:THR:HG22	1:B:579:MET:HE1	1.76	0.67
1:A:85:LEU:HA	1:A:88:ILE:HD12	1.78	0.66
1:B:423:LEU:H	1:B:423:LEU:HD12	1.60	0.65
1:B:568:LEU:C	1:B:568:LEU:HD23	2.21	0.64
1:A:197:GLU:HG3	1:B:197:GLU:CD	2.23	0.64
1:B:112:LYS:HG3	1:B:115:ARG:HH11	1.63	0.63
1:B:416:LYS:HD3	1:B:543:ASP:HB3	1.81	0.63
1:B:294:THR:HG23	1:B:365:THR:HA	1.81	0.62
1:A:517:THR:HG22	1:A:579:MET:HE1	1.81	0.62
1:A:450:LEU:HB2	1:A:451:PRO:HD3	1.81	0.62
1:A:56:GLU:O	1:A:59:VAL:HG22	2.01	0.60
1:B:369:PHE:O	1:B:373:HIS:HD2	1.85	0.60
1:A:549:GLU:H	1:A:549:GLU:CD	2.09	0.59
1:B:212:VAL:HG21	1:B:565:PRO:HG3	1.83	0.59
1:B:417:HIS:O	1:B:421:ILE:HG12	2.03	0.59
1:B:180:TYR:O	1:B:183:TYR:HB3	2.02	0.58
1:A:469:PRO:HB2	1:A:471:ASP:OD1	2.03	0.58
1:A:190:MET:HE1	1:A:202:TYR:CZ	2.39	0.57
1:B:212:VAL:HG11	1:B:565:PRO:HG2	1.85	0.57
1:A:524:GLN:HG2	1:A:583:PRO:HG2	1.86	0.57
1:B:477:TRP:CE3	1:B:500:PRO:HG3	2.38	0.57
1:A:137:ASN:HB3	1:A:140:GLU:HB2	1.85	0.57
1:A:134:ASN:OD1	1:A:135:PRO:HD2	2.04	0.57
1:B:406:GLU:HG3	1:B:518:ARG:HH11	1.70	0.57
1:B:323:MET:HA	1:B:323:MET:HE2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:VAL:HG22	1:B:193:ALA:HB1	1.89	0.55
1:B:293:VAL:HB	1:B:423:LEU:HB3	1.89	0.55
1:A:468:ILE:HG22	1:A:473:TRP:HD1	1.72	0.55
1:A:137:ASN:HB3	1:A:140:GLU:CB	2.38	0.54
1:B:557:MET:C	1:B:557:MET:SD	2.91	0.54
1:B:419:LYS:NZ	1:B:428:PHE:HB3	2.21	0.53
1:A:287:GLN:HA	1:A:287:GLN:OE1	2.08	0.53
1:B:523:PHE:CE2	1:B:584:LEU:HD12	2.44	0.53
1:B:589:GLU:HB3	1:B:590:PRO:HD3	1.91	0.53
1:A:132:VAL:O	1:A:142:LEU:N	2.38	0.52
1:B:227:GLU:HG2	1:B:454:TYR:OH	2.09	0.52
2:C:14:HIS:C	2:C:16:CYS:H	2.16	0.52
1:A:163:TRP:CD1	1:A:167:SER:HG	2.27	0.52
1:A:201:ASP:CG	1:A:219:ARG:HE	2.17	0.51
1:B:275:TRP:HB3	1:B:278:LEU:HD12	1.92	0.51
1:B:366:MET:O	1:B:369:PHE:HB3	2.11	0.51
1:B:390:PHE:CE1	2:D:4:ARG:HD3	2.45	0.51
1:A:240:LEU:HA	1:A:595:LEU:HD21	1.91	0.51
1:B:307:ILE:HG23	1:B:369:PHE:HD1	1.75	0.51
1:B:499:ASP:N	1:B:500:PRO:CD	2.74	0.51
1:A:142:LEU:HD12	1:A:163:TRP:HH2	1.77	0.50
1:A:477:TRP:CE3	1:A:500:PRO:HG3	2.45	0.50
1:B:371:THR:O	1:B:375:GLU:HG2	2.11	0.50
1:A:55:THR:O	1:A:59:VAL:HG13	2.11	0.50
1:A:455:MET:HE3	1:A:480:MET:HB2	1.94	0.50
1:B:38:ASP:O	1:B:42:GLN:HG2	2.12	0.50
1:A:177:ARG:N	1:A:178:PRO:HD2	2.26	0.50
1:B:85:LEU:HD22	1:B:94:LYS:HG3	1.93	0.50
1:B:315:PHE:HE2	1:B:376:MET:HB3	1.77	0.50
1:B:302:TRP:CG	1:B:306:ARG:HG2	2.47	0.50
1:B:285:PHE:CD2	1:B:433:GLU:HB3	2.47	0.49
1:A:86:GLN:HG3	1:A:87:GLU:N	2.27	0.49
1:B:414:THR:HG21	1:B:542:CYS:O	2.13	0.49
1:B:500:PRO:O	1:B:506:VAL:HG21	2.12	0.49
1:A:50:TYR:CE2	1:A:59:VAL:HG12	2.47	0.49
1:B:293:VAL:O	1:B:297:MET:HG3	2.13	0.49
1:A:190:MET:HE1	1:A:202:TYR:CE1	2.48	0.49
1:A:499:ASP:N	1:A:500:PRO:CD	2.76	0.49
1:A:526:GLN:HA	1:A:526:GLN:OE1	2.13	0.49
1:A:239:HIS:CE1	1:A:596:LYS:HG2	2.48	0.49
1:A:538:PRO:HB2	1:A:541:LYS:HD3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:ARG:NH2	1:B:495:GLU:O	2.45	0.49
1:B:26:LYS:HE2	1:B:93:VAL:HG11	1.95	0.48
1:B:90:ASN:HB2	1:B:93:VAL:HG22	1.94	0.48
1:B:168:TRP:CG	1:B:270:MET:HE3	2.49	0.48
1:A:431:ASP:O	1:A:435:GLU:HG2	2.14	0.48
1:B:279:TYR:CZ	1:B:283:VAL:HG23	2.49	0.48
1:A:428:PHE:CE2	1:A:430:GLU:HG2	2.49	0.48
1:B:192:ARG:HA	1:B:196:TYR:O	2.13	0.48
1:B:315:PHE:CE2	1:B:376:MET:HB3	2.48	0.48
1:B:307:ILE:CG2	1:B:369:PHE:HA	2.44	0.47
1:A:133:CYS:HA	1:A:140:GLU:O	2.13	0.47
1:B:474:MET:HE3	1:B:474:MET:HA	1.96	0.47
2:C:14:HIS:C	2:C:16:CYS:N	2.68	0.47
1:B:237:TYR:CE1	1:B:451:PRO:HG2	2.48	0.47
1:A:545:SER:O	1:A:546:ASN:HB2	2.15	0.47
1:A:284:PRO:HB3	1:A:594:TRP:CZ2	2.49	0.47
1:B:22:GLU:O	1:B:26:LYS:HD3	2.15	0.47
1:A:431:ASP:OD1	1:A:434:THR:HG23	2.14	0.46
1:B:50:TYR:CD2	1:B:50:TYR:C	2.94	0.46
1:A:240:LEU:HD22	1:A:447:VAL:HG22	1.98	0.46
1:A:126:ILE:HD11	1:A:176:LEU:CD2	2.46	0.46
1:A:249:MET:HG2	1:A:256:ILE:HG22	1.98	0.46
1:B:307:ILE:HG21	1:B:369:PHE:HA	1.97	0.46
1:B:327:PHE:HE2	1:B:358:ILE:HG13	1.81	0.46
1:B:295:ASP:OD1	1:B:295:ASP:N	2.49	0.46
1:A:133:CYS:HA	1:A:141:CYS:HA	1.97	0.46
1:B:245:ARG:NH2	1:B:258:PRO:O	2.48	0.46
1:B:305:GLN:O	1:B:309:LYS:HB2	2.16	0.46
1:A:26:LYS:HD2	1:A:93:VAL:HG11	1.98	0.45
1:B:145:GLU:HA	1:B:146:PRO:HA	1.84	0.45
1:A:197:GLU:HG3	1:B:197:GLU:HG3	1.98	0.45
1:B:133:CYS:HA	1:B:141:CYS:HA	1.99	0.45
1:B:112:LYS:CG	1:B:115:ARG:HH11	2.29	0.45
1:B:317:SER:O	1:B:546:ASN:HA	2.17	0.45
1:A:305:GLN:O	1:A:309:LYS:HB2	2.17	0.45
1:A:490:PRO:HA	1:A:612:PRO:HG2	1.98	0.45
1:B:456:LEU:HD22	1:B:512:PHE:CD2	2.52	0.44
1:B:415:PRO:O	1:B:416:LYS:C	2.61	0.44
1:B:92:THR:HG22	1:B:96:GLN:HE21	1.82	0.44
1:A:137:ASN:ND2	1:A:139:GLN:H	2.15	0.44
1:B:347:THR:HG22	1:B:349:TRP:NE1	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345:HIS:O	1:B:346:PRO:C	2.61	0.44
1:A:309:LYS:O	1:A:313:LYS:HG3	2.18	0.44
1:B:22:GLU:HG2	1:B:26:LYS:HE3	1.99	0.44
1:A:126:ILE:HD11	1:A:176:LEU:HD23	1.99	0.43
1:A:474:MET:HE1	1:A:499:ASP:HB2	2.00	0.43
1:B:351:LEU:HD12	1:B:351:LEU:N	2.33	0.43
1:A:25:ALA:HB1	1:A:97:LEU:HD11	2.00	0.43
1:B:91:LEU:O	1:B:95:LEU:HG	2.18	0.43
1:B:423:LEU:HD12	1:B:423:LEU:N	2.31	0.43
1:B:579:MET:HE2	1:B:579:MET:HB3	1.76	0.43
1:B:294:THR:CG2	1:B:365:THR:HA	2.47	0.43
1:B:511:SER:O	1:B:514:ARG:HD2	2.17	0.43
1:B:132:VAL:O	1:B:141:CYS:HA	2.19	0.43
1:B:557:MET:SD	1:B:557:MET:O	2.77	0.43
1:A:410:LEU:HD23	1:A:410:LEU:HA	1.90	0.43
1:B:499:ASP:N	1:B:500:PRO:HD2	2.34	0.43
1:B:55:THR:OG1	1:B:56:GLU:N	2.51	0.43
1:B:538:PRO:HB2	1:B:540:HIS:CD2	2.54	0.43
1:A:53:ASN:O	1:A:58:ASN:ND2	2.51	0.43
1:A:518:ARG:HD2	1:A:518:ARG:C	2.43	0.43
1:A:477:TRP:CD2	1:A:500:PRO:HG3	2.54	0.42
1:B:232:GLU:HB2	1:B:581:VAL:HG11	2.01	0.42
1:A:454:TYR:CZ	1:A:458:LYS:HD2	2.55	0.42
1:B:165:TRP:CH2	1:B:490:PRO:HD2	2.54	0.42
1:A:86:GLN:HG3	1:A:87:GLU:HG3	2.01	0.42
1:A:457:GLU:OE1	1:A:460:ARG:NH1	2.46	0.42
1:A:50:TYR:CD2	1:A:59:VAL:HG12	2.54	0.42
1:A:197:GLU:HG2	1:B:219:ARG:HH21	1.85	0.42
1:B:188:ASN:O	1:B:192:ARG:HG3	2.19	0.42
1:B:568:LEU:C	1:B:568:LEU:CD2	2.91	0.42
1:A:284:PRO:HB3	1:A:594:TRP:CH2	2.55	0.42
1:A:455:MET:HE1	1:A:477:TRP:CE3	2.55	0.42
1:B:51:ASN:HD21	2:D:11:MET:CE	2.33	0.42
1:B:547:SER:HB3	1:B:550:ALA:HB3	2.02	0.42
1:A:85:LEU:HD22	1:A:94:LYS:HG3	2.01	0.42
1:A:145:GLU:CD	1:A:146:PRO:HA	2.44	0.42
1:A:482:ARG:NH1	1:A:488:VAL:HG23	2.35	0.42
1:B:520:LEU:HD23	1:B:579:MET:HG3	2.03	0.41
1:A:554:LEU:O	1:A:555:PHE:C	2.62	0.41
1:B:309:LYS:HD2	1:B:328:TRP:CZ2	2.55	0.41
1:B:455:MET:HE3	1:B:480:MET:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:GLU:HG3	1:B:197:GLU:CG	2.49	0.41
1:B:177:ARG:HB3	1:B:178:PRO:HD3	2.01	0.41
1:B:397:ASN:OD1	1:B:397:ASN:C	2.63	0.41
1:B:302:TRP:CD2	1:B:306:ARG:HG2	2.56	0.41
1:A:589:GLU:HB3	1:A:590:PRO:HD3	2.01	0.41
1:B:401:HIS:CE1	2:D:8:HIS:CE1	3.09	0.41
1:A:476:LYS:HD2	1:A:476:LYS:HA	1.91	0.41
1:B:82:MET:O	1:B:82:MET:HG2	2.20	0.41
1:B:107:VAL:CG2	1:B:193:ALA:HB1	2.50	0.41
1:B:244:VAL:O	1:B:248:LEU:HG	2.20	0.41
1:B:366:MET:O	1:B:369:PHE:N	2.49	0.41
1:B:517:THR:HG22	1:B:579:MET:CE	2.49	0.41
1:B:555:PHE:O	1:B:559:ARG:HG2	2.21	0.41
1:A:440:LEU:C	1:A:440:LEU:HD13	2.46	0.41
1:B:137:ASN:ND2	1:B:140:GLU:HB2	2.36	0.41
1:A:168:TRP:CG	1:A:270:MET:HE3	2.55	0.40
1:A:170:SER:HA	1:A:174:LYS:HG3	2.04	0.40
1:B:306:ARG:HG3	1:B:307:ILE:N	2.36	0.40
1:A:470:LYS:HA	1:A:473:TRP:CD1	2.56	0.40
1:A:595:LEU:HD12	1:A:595:LEU:HA	1.82	0.40
1:A:147:GLY:O	1:A:150:GLU:HB3	2.22	0.40
1:B:234:LYS:HB3	1:B:235:PRO:HD3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	594/609 (98%)	571 (96%)	22 (4%)	1 (0%)	43	44
1	B	593/609 (97%)	560 (94%)	32 (5%)	1 (0%)	43	44
2	C	16/18 (89%)	14 (88%)	2 (12%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	16/18 (89%)	14 (88%)	2 (12%)	0	100	100
All	All	1219/1254 (97%)	1159 (95%)	58 (5%)	2 (0%)	43	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	LEU
1	B	537	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	314/538 (58%)	306 (98%)	8 (2%)	42	47
1	B	197/538 (37%)	185 (94%)	12 (6%)	17	14
2	C	14/14 (100%)	13 (93%)	1 (7%)	13	10
2	D	14/14 (100%)	13 (93%)	1 (7%)	13	10
All	All	539/1104 (49%)	517 (96%)	22 (4%)	27	27

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	SER
1	A	68	LYS
1	A	89	GLN
1	A	118	THR
1	A	280	SER
1	A	287	GLN
1	A	309	LYS
1	A	347	THR
1	B	394	ASN
1	B	401	HIS
1	B	447	VAL
1	B	449	THR

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Mol	Chain	Res	Type
1	B	465	LYS
1	B	482	ARG
1	B	483	GLU
1	B	531	GLN
1	B	558	LEU
1	B	560	LEU
1	B	573	VAL
1	B	584	LEU
2	C	7	CYS
2	D	7	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	76	GLN
1	A	86	GLN
1	A	101	GLN
1	A	137	ASN
1	A	139	GLN
1	A	149	ASN
1	A	195	HIS
1	A	228	HIS
1	B	417	HIS
1	B	540	HIS
2	C	5	GLN
2	C	14	HIS
2	D	14	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

2 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$
3	LFI	D	101	2	18,18,21	0.38	0	24,24,27	1.23	3 (12%)
3	LFI	C	101	2	18,18,21	0.72	0	24,24,27	1.46	2 (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
3	LFI	D	101	2	-	6/18/30/33	0/0/1/1
3	LFI	C	101	2	-	5/18/30/33	0/0/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	101	LFI	N3-C2-N2	-5.00	103.00	110.77
3	D	101	LFI	N3-C2-N2	-3.13	105.91	110.77
3	C	101	LFI	N2-C1-N1	-3.01	106.09	110.77
3	D	101	LFI	N2-C1-N1	-2.76	106.48	110.77
3	D	101	LFI	N3-C3-N1	-2.03	107.61	110.77

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	101	LFI	C8-C4-N1-C1
3	C	101	LFI	C8-C4-N1-C3
3	C	101	LFI	O2-C4-N1-C1

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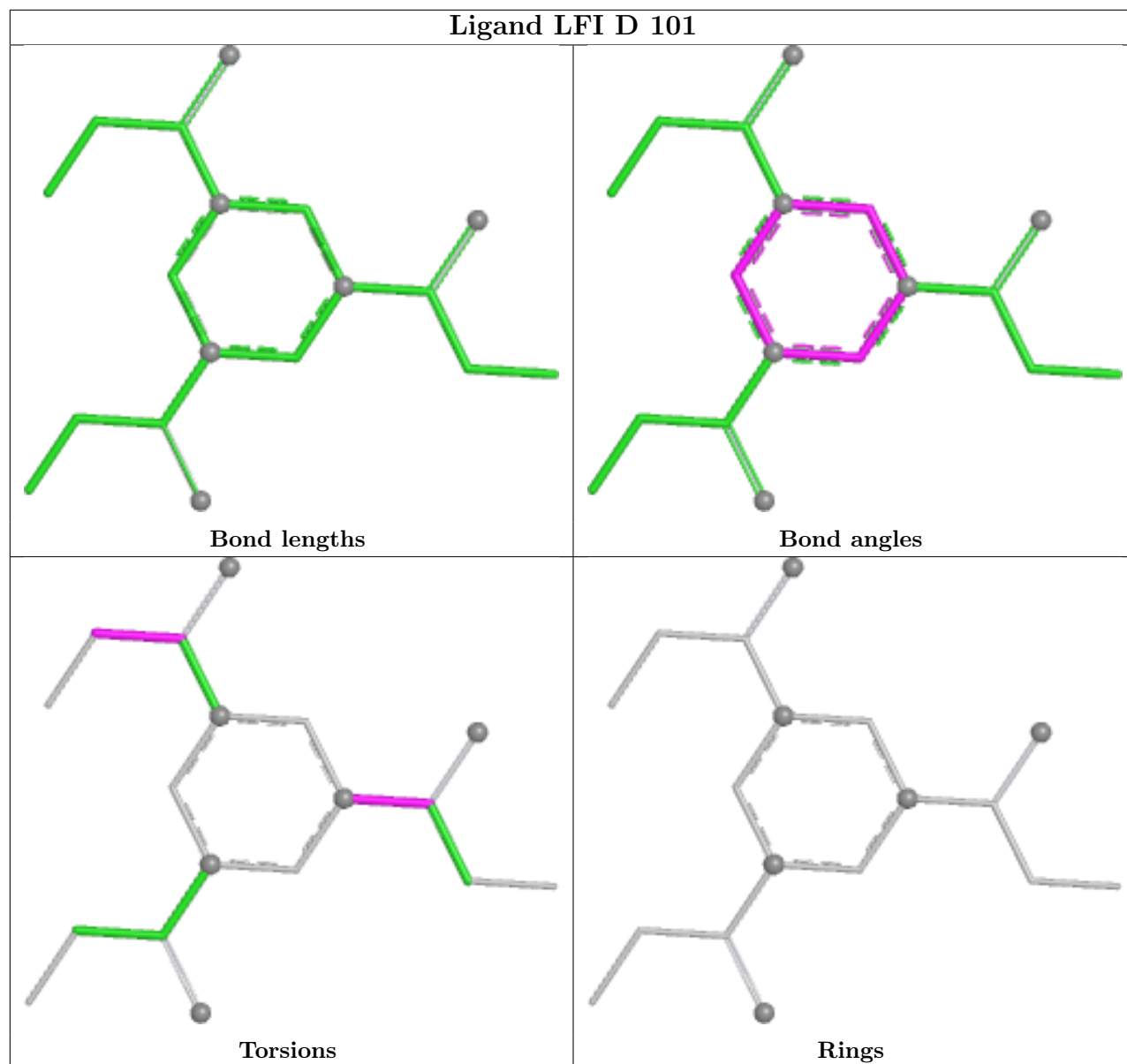
Mol	Chain	Res	Type	Atoms
3	C	101	LFI	O2-C4-N1-C3
3	D	101	LFI	C9-C5-N3-C3
3	D	101	LFI	O3-C5-N3-C3
3	D	101	LFI	O3-C5-N3-C2
3	D	101	LFI	N1-C4-C8-C10
3	D	101	LFI	O2-C4-C8-C10
3	D	101	LFI	C9-C5-N3-C2
3	C	101	LFI	O1-C6-C7-C12

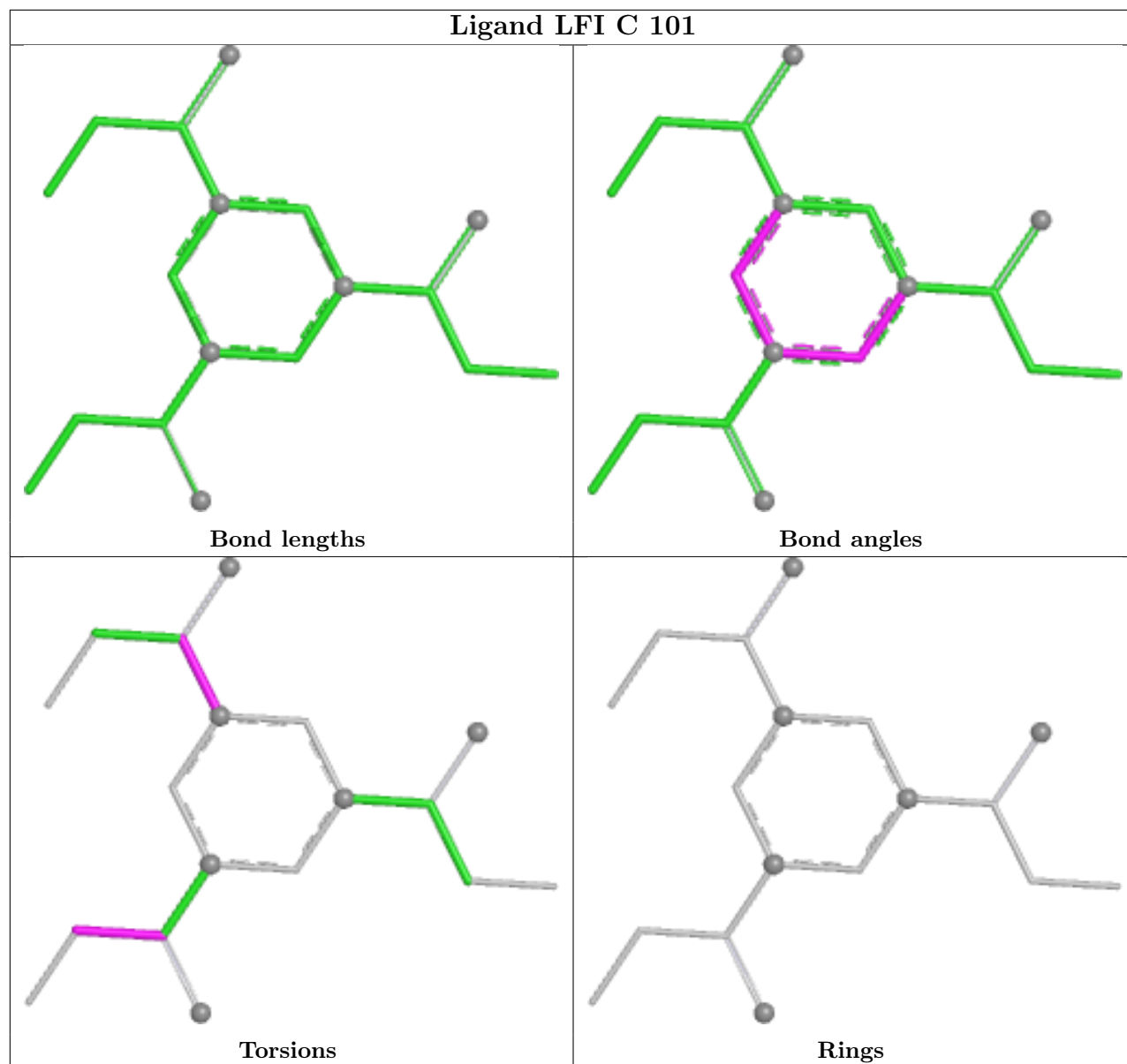
There are no ring outliers.

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.

Ligand LFI D 101





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	596/609 (97%)	0.43	27 (4%)	38 41	35, 53, 85, 158	0
1	B	595/609 (97%)	0.74	40 (6%)	24 26	37, 64, 108, 165	0
2	C	17/18 (94%)	0.75	2 (11%)	9 10	45, 59, 100, 124	0
2	D	17/18 (94%)	0.96	1 (5%)	28 30	49, 65, 90, 133	0
All	All	1225/1254 (97%)	0.59	70 (5%)	29 32	35, 59, 100, 165	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	108	LEU	5.0
1	B	107	VAL	4.4
1	B	105	SER	4.1
2	D	17	ALA	4.1
1	B	138	PRO	4.1
1	A	138	PRO	3.9
2	C	17	ALA	3.9
1	B	592	PHE	3.8
1	A	342	ALA	3.8
1	B	603	PHE	3.8
1	B	100	LEU	3.8
1	A	614	ALA	3.4
1	B	307	ILE	3.4
1	B	339	VAL	3.3
1	A	91	LEU	3.3
1	A	295	ASP	3.2
1	A	143	LEU	3.2
1	A	536	GLU	3.1
1	B	342	ALA	3.0
1	A	291	ILE	3.0
1	A	332	MET	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	432	ASN	2.9
1	B	343	VAL	2.9
1	B	428	PHE	2.8
1	A	213	ASP	2.8
1	A	126	ILE	2.8
1	A	142	LEU	2.8
1	A	343	VAL	2.8
1	A	573	VAL	2.7
1	A	156	LEU	2.7
1	B	423	LEU	2.7
1	B	215	TYR	2.7
1	A	339	VAL	2.7
1	B	291	ILE	2.7
1	B	302	TRP	2.6
1	B	413	ALA	2.6
1	A	21	ILE	2.5
1	B	259	ILE	2.5
1	A	334	THR	2.5
1	B	602	SER	2.5
1	B	103	ASN	2.5
1	A	135	PRO	2.5
1	B	308	PHE	2.5
1	A	289	PRO	2.4
1	B	400	PHE	2.3
1	B	212	VAL	2.3
1	A	336	PRO	2.3
2	C	15	LEU	2.3
1	B	239	HIS	2.3
1	A	88	ILE	2.3
1	B	21	ILE	2.2
1	B	121	ASN	2.2
1	A	211	GLY	2.2
1	B	243	TYR	2.2
1	B	20	THR	2.2
1	B	84	PRO	2.2
1	B	593	THR	2.2
1	B	594	TRP	2.2
1	A	412	ALA	2.1
1	A	224	GLU	2.1
1	B	91	LEU	2.1
1	B	289	PRO	2.1
1	B	364	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	126	ILE	2.1
1	B	596	LYS	2.1
1	B	23	GLU	2.0
1	B	85	LEU	2.0
1	A	132	VAL	2.0
1	A	390	PHE	2.0
1	B	304	ALA	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](#)

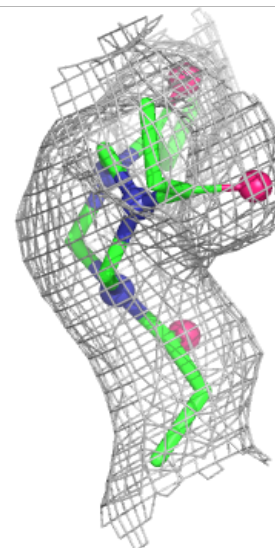
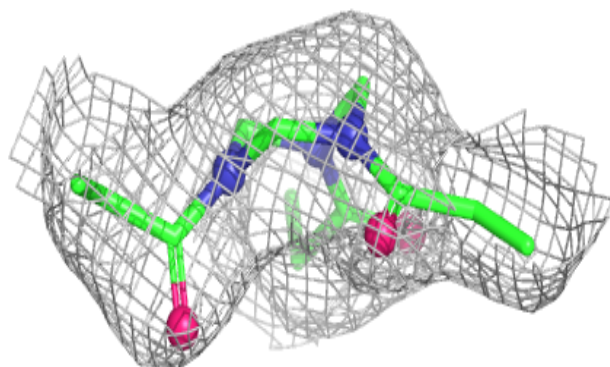
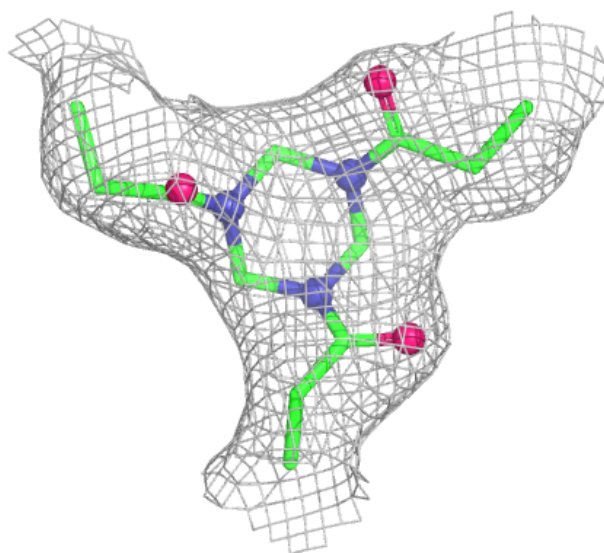
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	LFI	C	101	18/21	0.90	0.09	50,54,75,83	0
3	LFI	D	101	18/21	0.90	0.10	62,69,87,88	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.

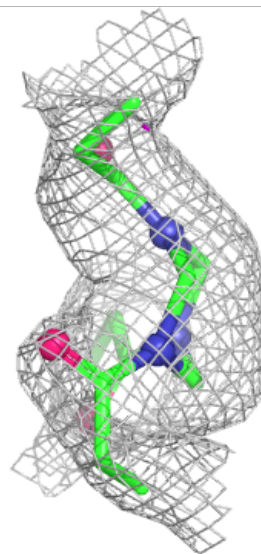
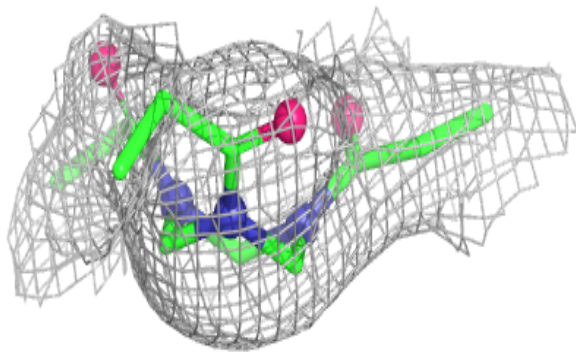
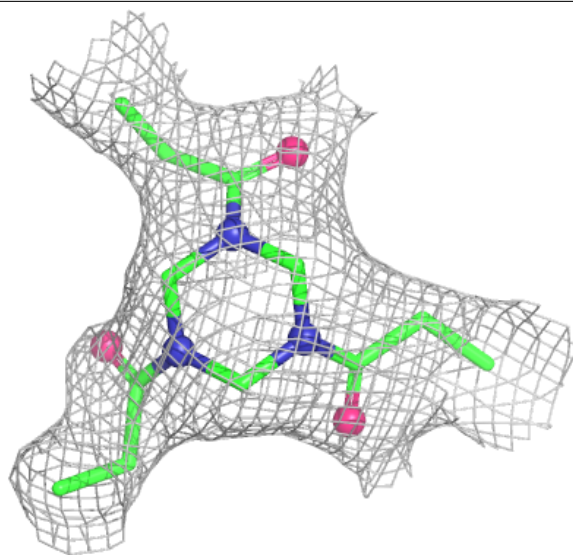
Electron density around LFI C 101:

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



Electron density around LFI D 101:

$2mF_o-DF_c$ (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.