



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:23 AM UTC

PDB ID : 4PJD / pdb_00004pjd
Title : Structure of human MR1-5-OP-RU in complex with human MAIT C-C10 TCR
Authors : Birkinshaw, R.W.; Rossjohn, J.
Deposited on : 2014-05-12
Resolution : 2.78 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

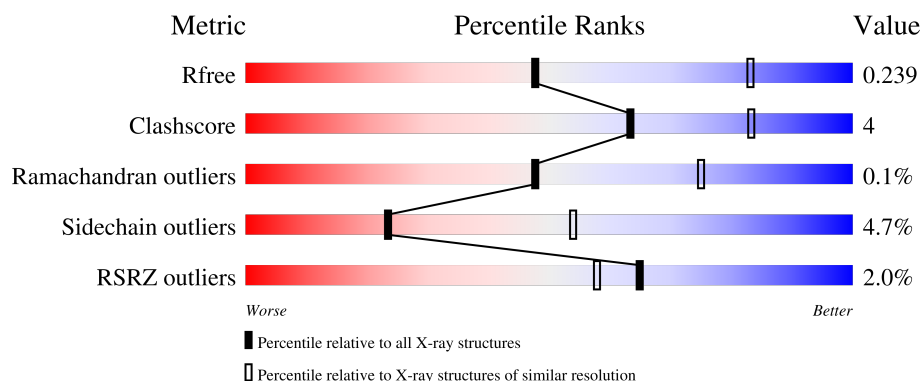
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	271	2% 80% 16% ..
1	C	271	% 79% 14% 6%
2	B	100	83% 12% ..
2	D	100	% 82% 12% ..
3	E	205	8% 80% 12% 8%

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Mol	Chain	Length	Quality of chain
3	G	205	85% 10% . .
4	F	245	2% 81% 13% . 5%
4	H	245	83% 13% . .

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12906 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 Major histocompatibility complex class I-related gene protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2147	1377	371	388	11			
1	C	255	Total	C	N	O	S	0	0	0
			2083	1331	363	378	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q95460
A	261	SER	CYS	engineered mutation	UNP Q95460
C	0	MET	-	initiating methionine	UNP Q95460
C	261	SER	CYS	engineered mutation	UNP Q95460

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	96	Total	C	N	O	S	0	0	0
			769	494	131	142	2			
2	D	96	Total	C	N	O	S	0	0	0
			753	484	129	138	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
D	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called TCR-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	188	Total	C	N	O	S	0	0	0
			1384	886	223	267	8			

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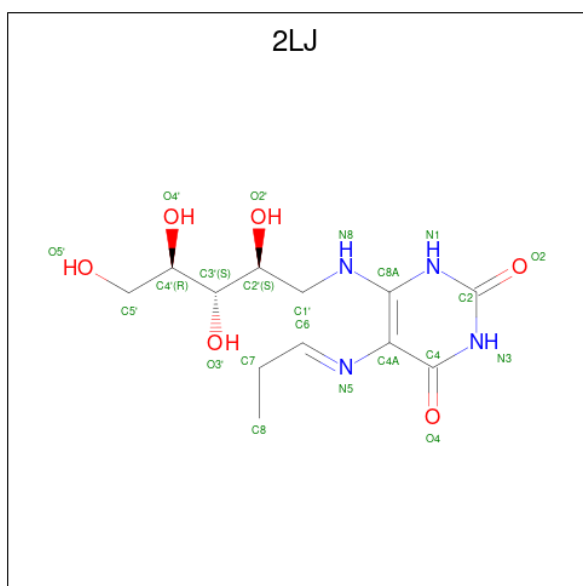
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	197	Total	C	N	O	S	0	0	0
			1495	949	240	297	9			

- Molecule 4 is a protein called TCR-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	F	233	Total	C	N	O	S	0	0	0
			1774	1120	305	340	9			
4	H	239	Total	C	N	O	S	0	0	0
			1821	1150	312	350	9			

- Molecule 5 is 1-deoxy-1-({2,6-dioxo-5-[(E)-propylideneamino]-1,2,3,6-tetrahydropyrimidin-4-yl}amino)-D-ribose (CCD ID: 2LJ) (formula: C₁₂H₂₀N₄O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			22	12	4	6		
5	C	1	Total	C	N	O	0	0
			22	12	4	6		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	C	O	0	0
			6	3	3		

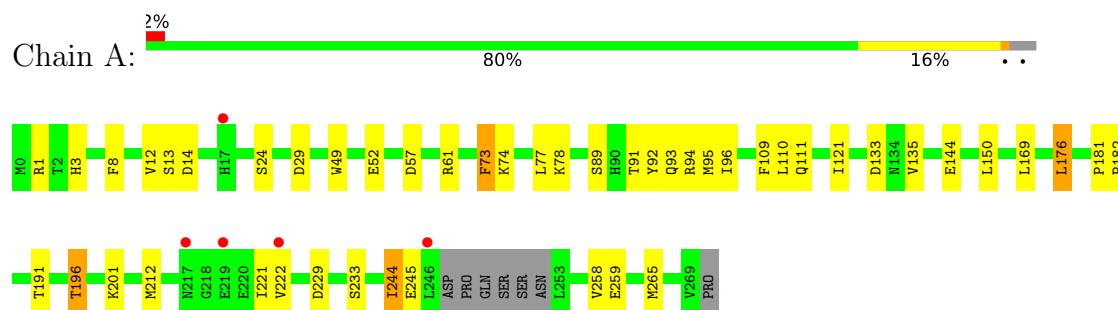
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	121	Total	O	0	0
			121	121		
7	B	61	Total	O	0	0
			61	61		
7	C	109	Total	O	0	0
			109	109		
7	D	33	Total	O	0	0
			33	33		
7	E	59	Total	O	0	0
			59	59		
7	F	87	Total	O	0	0
			87	87		
7	G	77	Total	O	0	0
			77	77		
7	H	83	Total	O	0	0
			83	83		

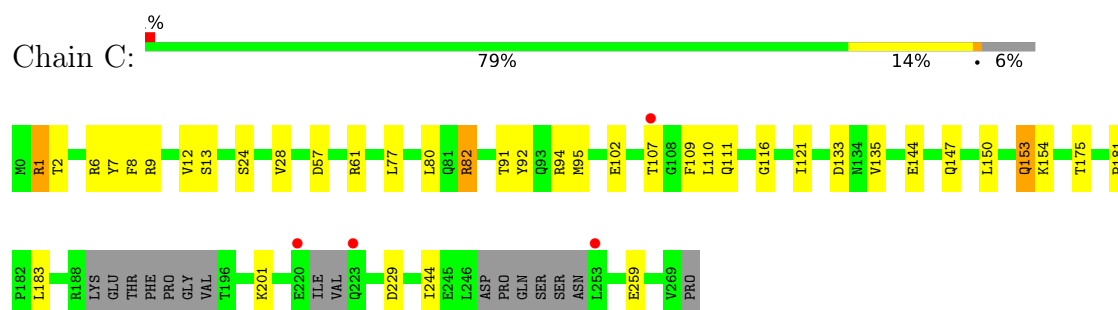
3 Residue-property plots [i](#)

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.

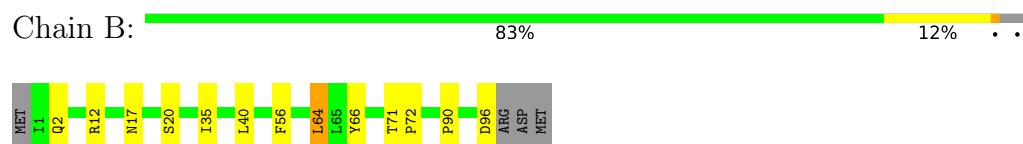
- Molecule 1: Major histocompatibility complex class I-related gene protein



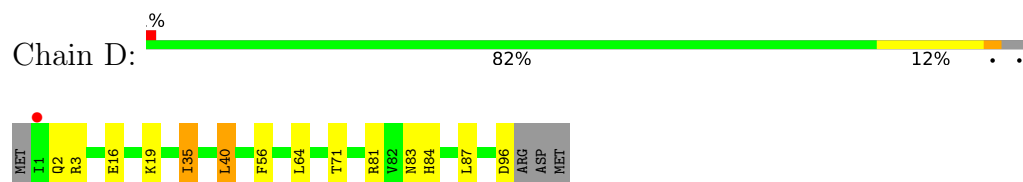
- Molecule 1: Major histocompatibility complex class I-related gene protein



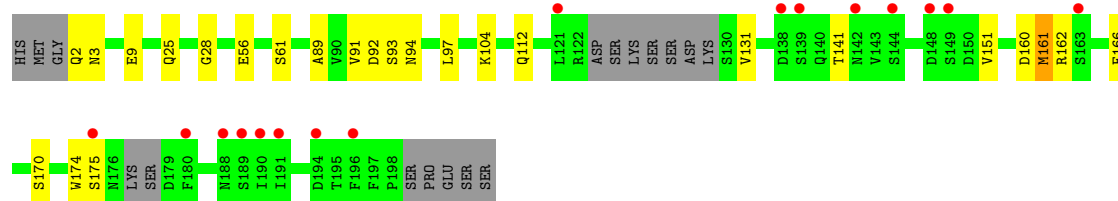
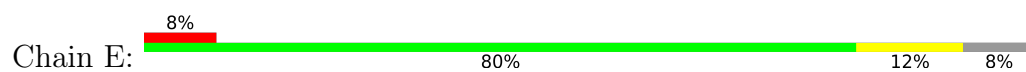
- Molecule 2: Beta-2-microglobulin



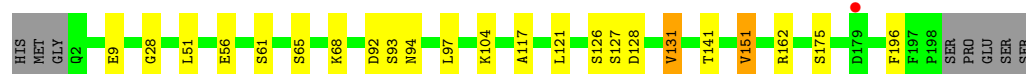
- Molecule 2: Beta-2-microglobulin



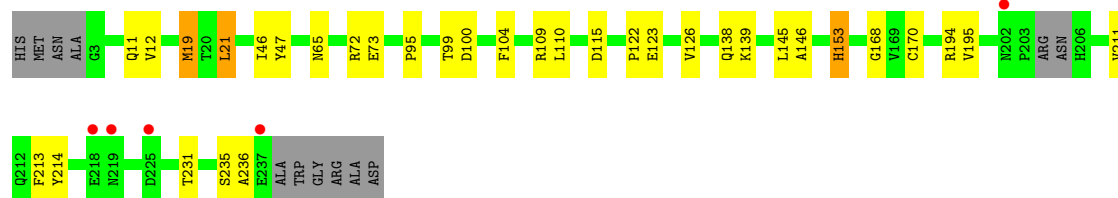
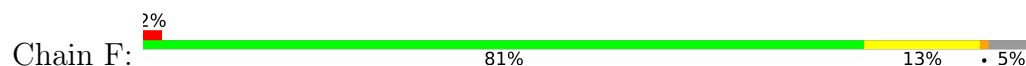
- Molecule 3: TCR-alpha



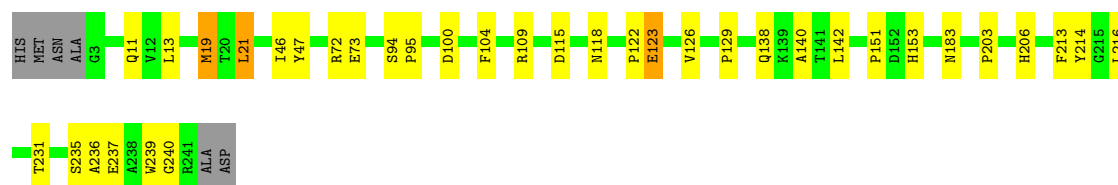
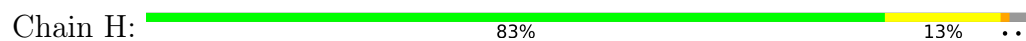
• Molecule 3: TCR-alpha



• Molecule 4: TCR-beta



• Molecule 4: TCR-beta



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	216.29Å 70.52Å 144.16Å 90.00° 104.50° 90.00°	Depositor
Resolution (Å)	40.10 – 2.78 40.10 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.10-2.78) 99.7 (40.10-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.77Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.168 , 0.226 0.178 , 0.239	Depositor DCC
R_{free} test set	2703 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	27.9	Xtriage
Anisotropy	1.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12906	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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: GOL, 2LJ

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	0.88	0/2212	1.30	12/3009 (0.4%)
1	C	0.86	0/2143	1.31	8/2907 (0.3%)
2	B	0.86	0/792	1.22	3/1080 (0.3%)
2	D	0.86	0/776	1.23	2/1061 (0.2%)
3	E	0.93	1/1416 (0.1%)	1.20	5/1932 (0.3%)
3	G	0.91	0/1529	1.31	10/2082 (0.5%)
4	F	0.82	0/1822	1.21	7/2491 (0.3%)
4	H	0.85	0/1872	1.24	6/2559 (0.2%)
All	All	0.87	1/12562 (0.0%)	1.26	53/17121 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	161	MET	SD-CE	-7.24	1.61	1.79

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	13	SER	N-CA-C	8.17	119.81	111.07
3	G	127	SER	CA-C-N	7.59	131.06	120.29
3	G	127	SER	C-N-CA	7.59	131.06	120.29
4	F	153	HIS	CA-CB-CG	7.15	120.95	113.80
3	G	128	ASP	CA-CB-CG	7.13	119.73	112.60
3	G	151	VAL	N-CA-CB	6.84	119.20	110.99
3	E	92	ASP	CA-CB-CG	6.53	119.13	112.60
1	C	28	VAL	N-CA-C	-6.50	97.74	107.37
2	D	71	THR	CB-CA-C	6.41	116.86	110.33
2	B	71	THR	CB-CA-C	6.28	116.74	110.33
3	G	162	ARG	CA-C-N	6.21	128.92	120.54
3	G	162	ARG	C-N-CA	6.21	128.92	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	128	ASP	CA-C-N	6.15	129.61	120.87
3	G	128	ASP	C-N-CA	6.15	129.61	120.87
1	A	229	ASP	CA-CB-CG	6.07	118.67	112.60
3	G	131	VAL	N-CA-C	6.02	116.54	108.11
1	C	135	VAL	N-CA-C	-5.80	105.08	110.53
4	F	65	ASN	CA-CB-CG	5.76	118.36	112.60
4	H	73	GLU	N-CA-C	5.71	118.55	109.24
1	C	28	VAL	N-CA-CB	5.70	118.85	111.67
4	H	115	ASP	CA-CB-CG	5.69	118.29	112.60
4	F	73	GLU	N-CA-C	5.69	118.52	109.24
4	H	183	ASN	CA-CB-CG	5.67	118.27	112.60
1	C	229	ASP	CA-CB-CG	5.66	118.26	112.60
1	A	14	ASP	CA-CB-CG	5.64	118.24	112.60
1	A	73	PHE	CA-C-N	5.60	127.79	120.28
1	A	73	PHE	C-N-CA	5.60	127.79	120.28
4	H	94	SER	N-CA-C	5.56	113.22	108.22
1	A	176	LEU	N-CA-C	5.49	117.34	111.36
3	G	92	ASP	CA-CB-CG	5.42	118.02	112.60
4	H	123	GLU	N-CA-C	-5.38	99.75	108.52
4	H	104	PHE	N-CA-C	5.35	118.45	109.95
2	B	96	ASP	CA-CB-CG	5.35	117.95	112.60
4	F	138	GLN	CB-CG-CD	5.32	121.65	112.60
1	C	133	ASP	CA-CB-CG	5.30	117.90	112.60
4	F	123	GLU	N-CA-C	-5.25	99.97	108.52
3	E	162	ARG	CA-C-N	5.22	127.59	120.54
3	E	162	ARG	C-N-CA	5.22	127.59	120.54
1	A	196	THR	CB-CA-C	5.20	118.32	109.80
4	F	115	ASP	CA-CB-CG	5.19	117.79	112.60
4	F	104	PHE	N-CA-C	5.15	118.13	109.95
2	B	56	PHE	N-CA-C	5.13	117.33	109.07
1	A	259	GLU	CB-CG-CD	5.11	121.29	112.60
1	A	135	VAL	N-CA-C	-5.11	105.86	110.82
3	E	3	ASN	CA-C-N	5.11	129.83	123.14
3	E	3	ASN	C-N-CA	5.11	129.83	123.14
1	C	181	PRO	N-CA-C	5.11	115.37	110.47
1	A	222	VAL	CA-C-N	5.08	130.25	121.86
1	A	222	VAL	C-N-CA	5.08	130.25	121.86
1	A	133	ASP	CA-CB-CG	5.07	117.67	112.60
1	A	181	PRO	N-CA-C	5.04	115.31	110.47
1	C	2	THR	CB-CA-C	5.02	117.80	110.26
2	D	56	PHE	N-CA-C	5.00	117.12	109.07

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	2147	0	2023	20	0
1	C	2083	0	1940	22	0
2	B	769	0	715	5	0
2	D	753	0	682	5	0
3	E	1384	0	1216	16	0
3	G	1495	0	1369	6	0
4	F	1774	0	1632	17	0
4	H	1821	0	1674	14	0
5	A	22	0	18	1	0
5	C	22	0	18	5	0
6	E	6	0	8	0	0
7	A	121	0	0	1	0
7	B	61	0	0	1	0
7	C	109	0	0	2	0
7	D	33	0	0	0	0
7	E	59	0	0	0	0
7	F	87	0	0	0	0
7	G	77	0	0	0	0
7	H	83	0	0	1	0
All	All	12906	0	11295	95	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 (95) 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)
3:E:161:MET:HE3	3:E:166:PHE:HD2	1.39	0.88
3:E:161:MET:HE3	3:E:166:PHE:CD2	2.17	0.78
1:C:107:THR:N	1:C:107:THR:C	2.43	0.77
1:A:96:ILE:HD12	1:A:110:LEU:HD13	1.77	0.65
1:C:110:LEU:HD22	1:C:153:GLN:HG2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:THR:C	1:C:107:THR:CB	2.71	0.63
1:C:12:VAL:HG22	1:C:91:THR:HG22	1.81	0.63
3:E:161:MET:CE	4:F:194:ARG:HD3	2.28	0.63
3:E:161:MET:HE1	4:F:139:LYS:HD3	1.80	0.63
3:E:2:GLN:HB3	3:E:25:GLN:O	2.00	0.61
3:E:28:GLY:HA3	3:E:93:SER:HB3	1.85	0.58
3:E:161:MET:HE2	4:F:194:ARG:HD3	1.84	0.58
4:H:129:PRO:HD3	4:H:142:LEU:HG	1.86	0.58
4:F:126:VAL:HG23	4:F:236:ALA:HB3	1.84	0.57
4:H:126:VAL:HG23	4:H:236:ALA:HB3	1.84	0.57
1:C:144:GLU:HA	1:C:150:LEU:HD11	1.86	0.56
1:A:182:PRO:HG2	1:A:265:MET:HE1	1.87	0.56
1:C:147:GLN:NE2	1:C:150:LEU:HD12	2.22	0.55
1:C:1:ARG:HD3	7:C:506:HOH:O	2.07	0.55
4:F:12:VAL:HG22	4:F:153:HIS:CE1	2.42	0.55
1:A:144:GLU:HA	1:A:150:LEU:HD11	1.88	0.55
1:C:107:THR:N	1:C:107:THR:CB	2.70	0.54
5:A:301:2LJ:O4	5:A:301:2LJ:H1	2.07	0.54
4:H:138:GLN:HG3	7:H:343:HOH:O	2.07	0.54
3:E:161:MET:CE	3:E:166:PHE:CD2	2.91	0.53
3:E:9:GLU:HG3	3:E:104:LYS:HB3	1.90	0.53
4:H:151:PRO:HG2	4:H:153:HIS:CD2	2.44	0.53
1:A:94:ARG:HD3	7:A:441:HOH:O	2.09	0.52
3:G:9:GLU:HG3	3:G:104:LYS:HB3	1.92	0.52
4:H:11:GLN:HG2	4:H:19:MET:SD	2.49	0.52
2:B:64:LEU:HD22	2:B:66:TYR:HE1	1.74	0.52
4:H:129:PRO:HG3	4:H:140:ALA:HB1	1.93	0.51
4:F:11:GLN:HG2	4:F:19:MET:SD	2.51	0.51
1:A:12:VAL:HG22	1:A:91:THR:HG22	1.93	0.51
1:C:154:LYS:HD3	3:G:51:LEU:HD11	1.93	0.50
1:A:212:MET:HG2	1:A:258:VAL:HG22	1.94	0.50
3:E:160:ASP:O	4:F:168:GLY:N	2.42	0.50
4:H:153:HIS:HB3	4:H:214:TYR:HB2	1.93	0.50
1:A:95:MET:HE3	1:A:109:PHE:CD1	2.46	0.50
1:A:169:LEU:HD23	1:A:176:LEU:HD13	1.93	0.50
3:E:151:VAL:HA	3:E:175:SER:HB2	1.94	0.49
4:F:214:TYR:HA	4:F:231:THR:HG23	1.95	0.49
3:E:89:ALA:HB1	3:E:97:LEU:HD22	1.95	0.48
2:D:16:GLU:HB3	2:D:19:LYS:HG3	1.95	0.48
5:C:301:2LJ:O4	5:C:301:2LJ:H1	2.14	0.48
2:B:17:ASN:HA	2:B:72:PRO:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:153:HIS:HB3	4:F:214:TYR:HB2	1.96	0.47
2:D:40:LEU:HD11	2:D:81:ARG:HB2	1.97	0.47
4:H:206:HIS:HD2	4:H:239:TRP:CE2	2.32	0.47
1:A:233:SER:HB3	2:B:12:ARG:HG3	1.97	0.47
1:A:77:LEU:HD13	1:A:92:TYR:HB2	1.95	0.47
1:C:94:ARG:HH12	5:C:301:2LJ:H14	1.80	0.46
4:H:214:TYR:HA	4:H:231:THR:HG23	1.98	0.46
1:A:111:GLN:HG2	1:A:121:ILE:HG23	1.97	0.46
1:C:94:ARG:HD3	7:C:483:HOH:O	2.16	0.46
4:F:122:PRO:HD3	4:F:213:PHE:HB2	1.97	0.46
4:F:12:VAL:HG22	4:F:153:HIS:HE1	1.81	0.46
1:C:77:LEU:HD13	1:C:92:TYR:HB2	1.98	0.45
1:A:245:GLU:HG3	1:C:82:ARG:HH11	1.81	0.45
2:D:35:ILE:HD12	2:D:84:HIS:HD2	1.81	0.45
4:F:95:PRO:HD2	4:F:100:ASP:HB3	1.99	0.45
1:A:61:ARG:HD3	3:E:94:ASN:HB3	1.99	0.45
4:F:46:ILE:HG22	4:F:47:TYR:HD2	1.81	0.45
2:B:90:PRO:HG2	7:B:142:HOH:O	2.16	0.44
1:A:74:LYS:O	1:A:78:LYS:HG3	2.18	0.44
4:H:95:PRO:HD2	4:H:100:ASP:HB3	2.00	0.44
1:A:49:TRP:HA	1:A:52:GLU:HG2	2.00	0.43
1:C:9:ARG:HH22	5:C:301:2LJ:H16	1.83	0.43
1:C:7:TYR:OH	5:C:301:2LJ:H4	2.19	0.43
4:F:146:ALA:HB2	4:F:211:VAL:HG21	2.00	0.43
1:C:116:GLY:O	2:D:3:ARG:NH2	2.51	0.43
1:A:8:PHE:O	1:A:24:SER:HA	2.18	0.43
1:A:3:HIS:HA	1:A:29:ASP:OD1	2.19	0.43
1:C:94:ARG:HH22	5:C:301:2LJ:H17	1.84	0.43
1:C:111:GLN:HG2	1:C:121:ILE:HG23	2.01	0.43
3:E:161:MET:HE1	4:F:194:ARG:HD3	2.00	0.42
1:A:221:ILE:HD13	1:A:244:ILE:HG21	2.01	0.42
3:E:131:VAL:HG22	3:E:174:TRP:HB3	2.01	0.42
1:C:8:PHE:O	1:C:24:SER:HA	2.19	0.42
4:F:21:LEU:HD23	4:F:21:LEU:N	2.35	0.42
4:H:46:ILE:HG22	4:H:47:TYR:HD2	1.83	0.42
3:G:65:SER:OG	3:G:68:LYS:HB2	2.20	0.42
1:A:13:SER:O	1:A:89:SER:OG	2.36	0.41
3:G:117:ALA:HB2	3:G:196:PHE:HB3	2.02	0.41
2:D:83:ASN:HA	2:D:87:LEU:HD12	2.02	0.41
3:G:28:GLY:HA3	3:G:93:SER:HB3	2.03	0.41
1:C:95:MET:HE3	1:C:109:PHE:CD1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:ARG:HD3	3:G:94:ASN:HB3	2.02	0.41
4:H:21:LEU:N	4:H:21:LEU:HD23	2.36	0.41
2:B:64:LEU:HD22	2:B:66:TYR:CE1	2.56	0.40
3:E:91:VAL:HG11	4:F:99:THR:HG22	2.03	0.40
4:H:122:PRO:HD3	4:H:213:PHE:HB2	2.02	0.40
1:A:57:ASP:O	1:A:61:ARG:HG3	2.21	0.40
1:C:57:ASP:O	1:C:61:ARG:HG3	2.21	0.40
4:H:203:PRO:HA	4:H:240:GLY:O	2.22	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	260/271 (96%)	252 (97%)	7 (3%)	1 (0%)	30	57
1	C	246/271 (91%)	241 (98%)	5 (2%)	0	100	100
2	B	94/100 (94%)	92 (98%)	2 (2%)	0	100	100
2	D	94/100 (94%)	92 (98%)	2 (2%)	0	100	100
3	E	182/205 (89%)	175 (96%)	6 (3%)	1 (0%)	24	52
3	G	195/205 (95%)	190 (97%)	5 (3%)	0	100	100
4	F	229/245 (94%)	221 (96%)	8 (4%)	0	100	100
4	H	237/245 (97%)	230 (97%)	7 (3%)	0	100	100
All	All	1537/1642 (94%)	1493 (97%)	42 (3%)	2 (0%)	48	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	THR
3	E	112	GLN

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	224/241 (93%)	218 (97%)	6 (3%)	39	71
1	C	216/241 (90%)	205 (95%)	11 (5%)	21	50
2	B	83/95 (87%)	78 (94%)	5 (6%)	17	43
2	D	78/95 (82%)	73 (94%)	5 (6%)	16	41
3	E	133/182 (73%)	129 (97%)	4 (3%)	36	69
3	G	158/182 (87%)	149 (94%)	9 (6%)	18	46
4	F	186/212 (88%)	177 (95%)	9 (5%)	23	53
4	H	190/212 (90%)	180 (95%)	10 (5%)	20	49
All	All	1268/1460 (87%)	1209 (95%)	59 (5%)	23	54

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ARG
1	A	73	PHE
1	A	93	GLN
1	A	196	THR
1	A	201	LYS
1	A	244	ILE
2	B	2	GLN
2	B	20	SER
2	B	35	ILE
2	B	40	LEU
2	B	64	LEU
1	C	1	ARG
1	C	6	ARG
1	C	80	LEU
1	C	82	ARG
1	C	102	GLU
1	C	153	GLN
1	C	175	THR
1	C	183	LEU

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Mol	Chain	Res	Type
1	C	201	LYS
1	C	244	ILE
1	C	259	GLU
2	D	2	GLN
2	D	35	ILE
2	D	40	LEU
2	D	64	LEU
2	D	96	ASP
3	E	56	GLU
3	E	61	SER
3	E	141	THR
3	E	170	SER
4	F	19	MET
4	F	21	LEU
4	F	72	ARG
4	F	109	ARG
4	F	110	LEU
4	F	145	LEU
4	F	170	CYS
4	F	195	VAL
4	F	235	SER
3	G	56	GLU
3	G	61	SER
3	G	97	LEU
3	G	121	LEU
3	G	126	SER
3	G	131	VAL
3	G	141	THR
3	G	151	VAL
3	G	175	SER
4	H	13	LEU
4	H	19	MET
4	H	21	LEU
4	H	72	ARG
4	H	109	ARG
4	H	118	ASN
4	H	123	GLU
4	H	216	LEU
4	H	235	SER
4	H	237	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	137	HIS
1	A	148	HIS
2	B	31	HIS
1	C	111	GLN
1	C	134	ASN
1	C	147	GLN
1	C	148	HIS
1	C	177	GLN
1	C	203	HIS
1	C	239	GLN
3	E	19	GLN
3	E	21	ASN
3	E	25	GLN
3	E	38	HIS
3	E	96	GLN
4	F	22	GLN
4	F	102	GLN
3	G	19	GLN
3	G	21	ASN
3	G	38	HIS
3	G	96	GLN
4	H	138	GLN
4	H	153	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 ⓘ

3 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$
5	2LJ	C	301	1	22,22,22	1.50	3 (13%)	24,29,29	1.62	4 (16%)
6	GOL	E	301	-	5,5,5	0.10	0	5,5,5	0.17	0
5	2LJ	A	301	1	22,22,22	1.74	3 (13%)	24,29,29	1.84	8 (33%)

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
5	2LJ	C	301	1	-	11/18/19/19	0/1/1/1
6	GOL	E	301	-	-	0/4/4/4	-
5	2LJ	A	301	1	-	8/18/19/19	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	301	2LJ	C8A-N8	6.22	1.42	1.32
5	C	301	2LJ	C8A-N8	5.63	1.41	1.32
5	A	301	2LJ	C1'-C2'	3.95	1.57	1.52
5	C	301	2LJ	C7-C6	2.52	1.51	1.49
5	A	301	2LJ	C5'-C4'	2.51	1.58	1.52
5	C	301	2LJ	C1'-C2'	2.45	1.55	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	301	2LJ	C4'-C3'-C2'	5.13	122.10	113.57
5	C	301	2LJ	O4'-C4'-C3'	3.82	118.19	109.25
5	C	301	2LJ	C4'-C3'-C2'	3.45	119.31	113.57
5	C	301	2LJ	C2'-C1'-N8	3.42	121.02	111.53
5	A	301	2LJ	O5'-C5'-C4'	3.10	117.67	111.16
5	A	301	2LJ	C6-N5-C4A	2.58	124.84	121.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	301	2LJ	C1'-N8-C8A	-2.46	121.49	126.69
5	A	301	2LJ	C2'-C1'-N8	2.44	118.29	111.53
5	A	301	2LJ	O4'-C4'-C3'	2.34	114.73	109.25
5	A	301	2LJ	O4'-C4'-C5'	2.20	114.03	109.03
5	C	301	2LJ	O4'-C4'-C5'	2.08	113.76	109.03
5	A	301	2LJ	O2'-C2'-C1'	-2.07	105.59	109.67

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	301	2LJ	C2'-C3'-C4'-O4'
5	A	301	2LJ	O3'-C3'-C4'-O4'
5	C	301	2LJ	C1'-C2'-C3'-O3'
5	C	301	2LJ	C1'-C2'-C3'-C4'
5	C	301	2LJ	C2'-C3'-C4'-O4'
5	C	301	2LJ	O3'-C3'-C4'-O4'
5	C	301	2LJ	C3'-C4'-C5'-O5'
5	A	301	2LJ	O3'-C3'-C4'-C5'
5	A	301	2LJ	C2'-C3'-C4'-C5'
5	C	301	2LJ	O2'-C2'-C3'-C4'
5	C	301	2LJ	O3'-C3'-C4'-C5'
5	C	301	2LJ	O2'-C2'-C3'-O3'
5	C	301	2LJ	C2'-C3'-C4'-C5'
5	A	301	2LJ	O2'-C2'-C3'-O3'
5	A	301	2LJ	N8-C1'-C2'-C3'
5	C	301	2LJ	N8-C1'-C2'-C3'
5	C	301	2LJ	O4'-C4'-C5'-O5'
5	A	301	2LJ	C1'-C2'-C3'-O3'
5	A	301	2LJ	O2'-C2'-C3'-C4'

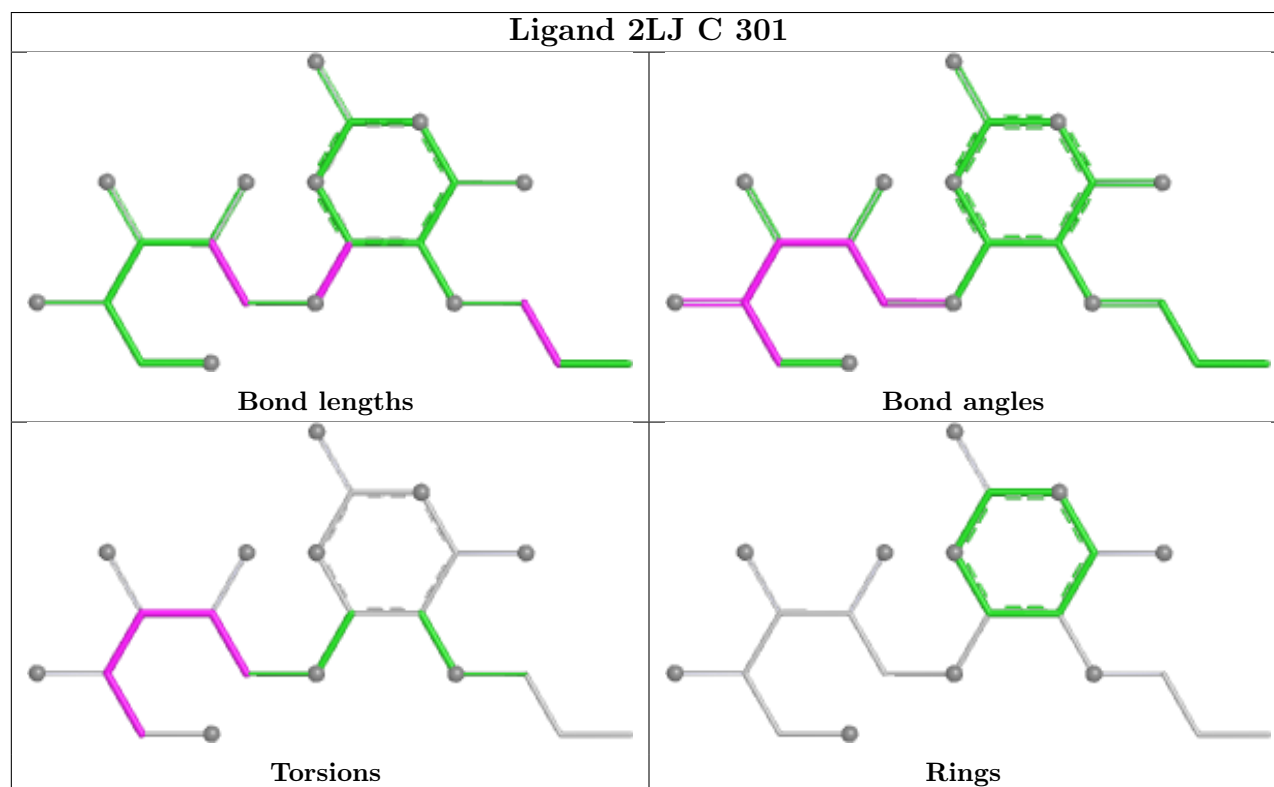
There are no ring outliers.

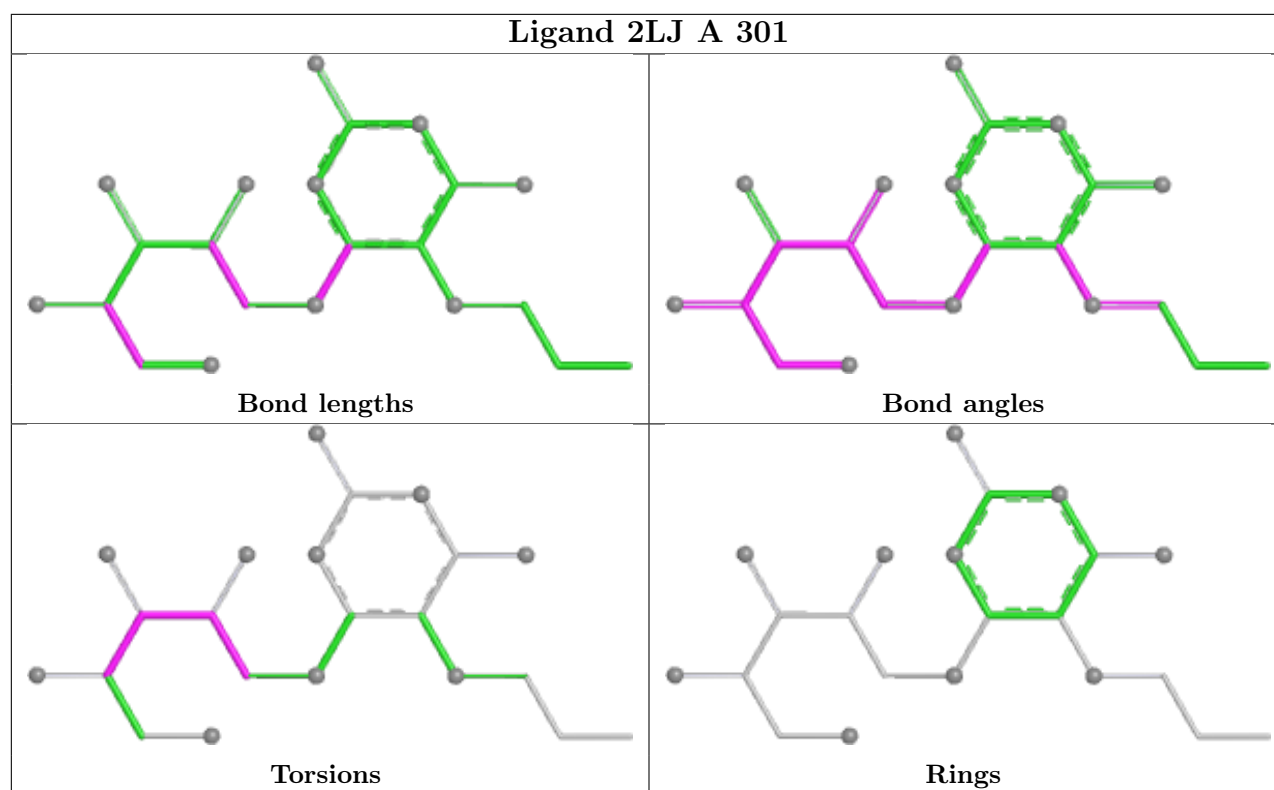
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	301	2LJ	5	0
5	A	301	2LJ	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 ⓘ

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	264/271 (97%)	-0.40	5 (1%) 66 59	10, 24, 48, 69	0
1	C	255/271 (94%)	-0.23	4 (1%) 70 63	13, 28, 56, 82	0
2	B	96/100 (96%)	-0.33	0 100 100	13, 30, 54, 64	0
2	D	96/100 (96%)	0.21	1 (1%) 79 74	18, 45, 75, 82	0
3	E	188/205 (91%)	0.22	16 (8%) 16 13	9, 36, 75, 93	0
3	G	197/205 (96%)	-0.31	1 (0%) 87 83	13, 26, 50, 80	0
4	F	233/245 (95%)	-0.07	5 (2%) 63 56	15, 34, 73, 90	0
4	H	239/245 (97%)	-0.39	0 100 100	12, 26, 46, 71	0
All	All	1568/1642 (95%)	-0.19	32 (2%) 65 58	9, 29, 66, 93	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	223	GLN	4.3
1	C	220	GLU	3.7
3	E	163	SER	3.3
3	E	148	ASP	3.2
1	C	107	THR	3.2
1	A	222	VAL	3.2
3	E	188	ASN	3.0
3	E	139	SER	2.9
3	E	149	SER	2.9
4	F	218	GLU	2.9
3	E	121	LEU	2.9
3	E	142	ASN	2.8
4	F	219	ASN	2.8
3	E	191	ILE	2.8
3	E	190	ILE	2.6
4	F	225	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
3	E	144	SER	2.4
2	D	1	ILE	2.4
1	A	17	HIS	2.4
1	C	253	LEU	2.3
1	A	246	LEU	2.3
3	E	180	PHE	2.3
1	A	217	ASN	2.2
3	E	175	SER	2.2
3	G	179	ASP	2.1
3	E	138	ASP	2.1
3	E	196	PHE	2.1
3	E	194	ASP	2.1
1	A	219	GLU	2.0
4	F	237	GLU	2.0
3	E	189	SER	2.0
4	F	202	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 [i](#)

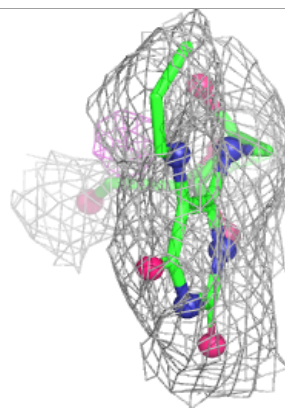
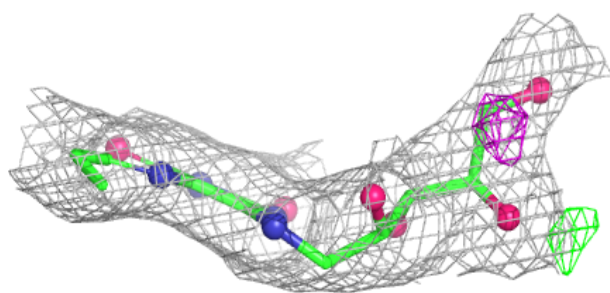
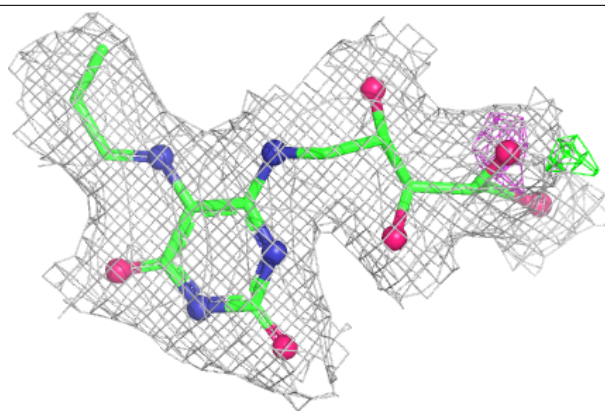
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
6	GOL	E	301	6/6	0.85	0.17	56,59,61,62	0
5	2LJ	A	301	22/22	0.95	0.08	7,12,18,20	0
5	2LJ	C	301	22/22	0.96	0.07	8,14,23,28	0

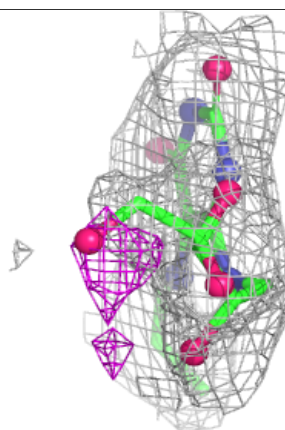
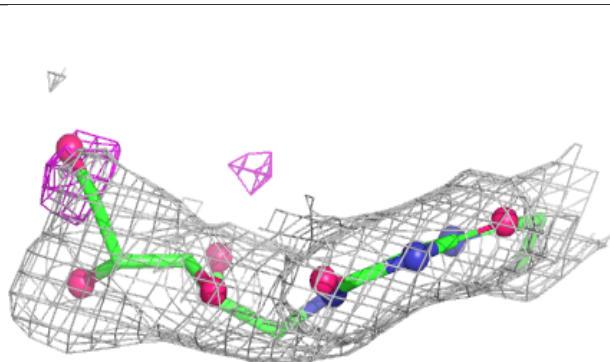
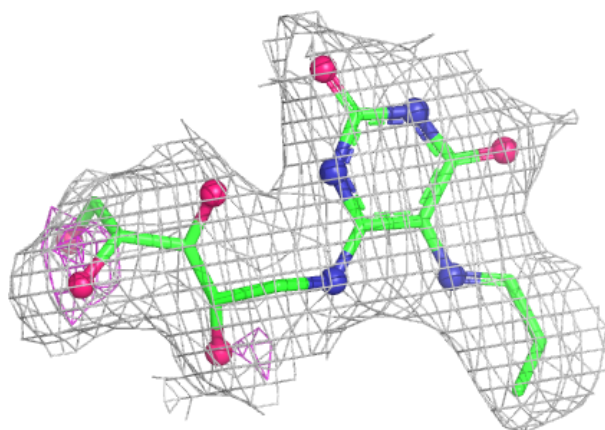
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 2LJ A 301:

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

**Electron density around 2LJ C 301:**

$2mF_o - DF_c$ (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.