



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:11 AM UTC

PDB ID : 6WTT / pdb_00006wtt
Title : Crystals Structure of the SARS-CoV-2 (COVID-19) main protease with inhibitor GC-376
Authors : Sacco, M.; Ma, C.; Chen, Y.; Wang, J.
Deposited on : 2020-05-03
Resolution : 2.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

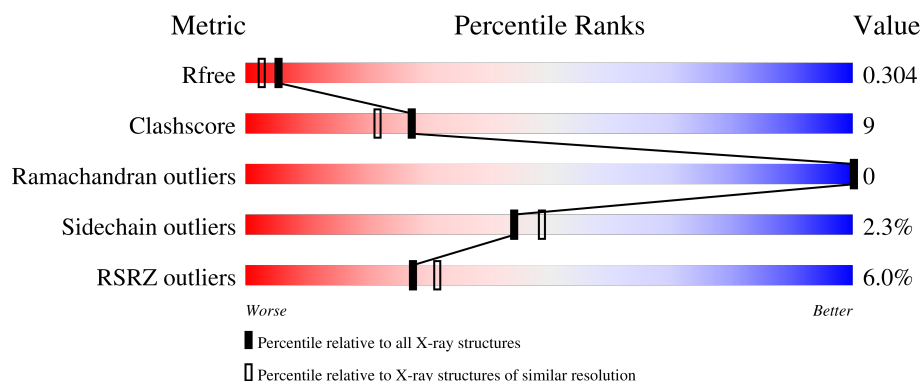
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	310	 2% 82% 15% ..
1	B	310	 10% 77% 19% ..
1	C	310	 6% 73% 24% ..

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7430 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.

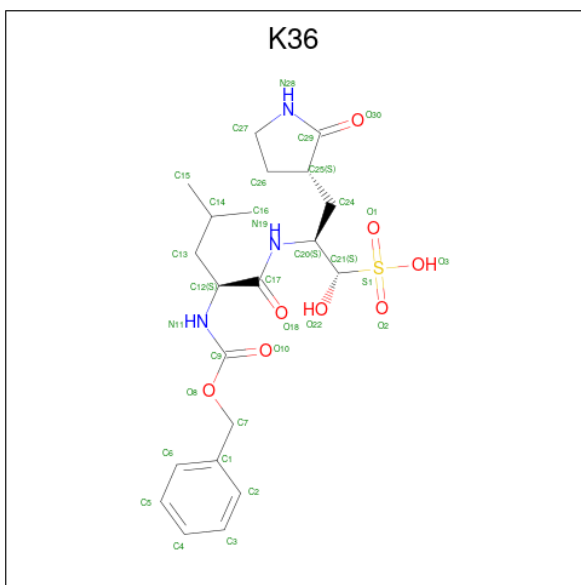
- Molecule 1 is a protein called 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2326	1470	397	437	22			
1	B	302	Total	C	N	O	S	0	0	0
			2333	1476	397	438	22			
1	C	302	Total	C	N	O	S	0	0	0
			2324	1468	397	437	22			

There are 18 discrepancies between the modelled and reference sequences:

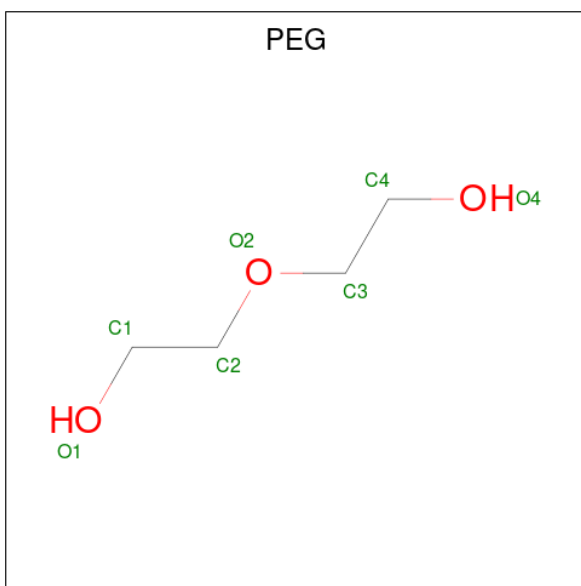
Chain	Residue	Modelled	Actual	Comment	Reference
A	305	HIS	-	expression tag	UNP P0DTD1
A	306	HIS	-	expression tag	UNP P0DTD1
A	307	HIS	-	expression tag	UNP P0DTD1
A	308	HIS	-	expression tag	UNP P0DTD1
A	309	HIS	-	expression tag	UNP P0DTD1
A	310	HIS	-	expression tag	UNP P0DTD1
B	305	HIS	-	expression tag	UNP P0DTD1
B	306	HIS	-	expression tag	UNP P0DTD1
B	307	HIS	-	expression tag	UNP P0DTD1
B	308	HIS	-	expression tag	UNP P0DTD1
B	309	HIS	-	expression tag	UNP P0DTD1
B	310	HIS	-	expression tag	UNP P0DTD1
C	305	HIS	-	expression tag	UNP P0DTD1
C	306	HIS	-	expression tag	UNP P0DTD1
C	307	HIS	-	expression tag	UNP P0DTD1
C	308	HIS	-	expression tag	UNP P0DTD1
C	309	HIS	-	expression tag	UNP P0DTD1
C	310	HIS	-	expression tag	UNP P0DTD1

- Molecule 2 is (1S,2S)-2-({N-[(benzyloxy)carbonyl]-L-leucyl}amino)-1-hydroxy-3-[(3S)-2-oxo pyrrolidin-3-yl]propane-1-sulfonic acid (CCD ID: K36) (formula: C₂₁H₃₁N₃O₈S) (labeled as "Ligand of Interest" by depositor).



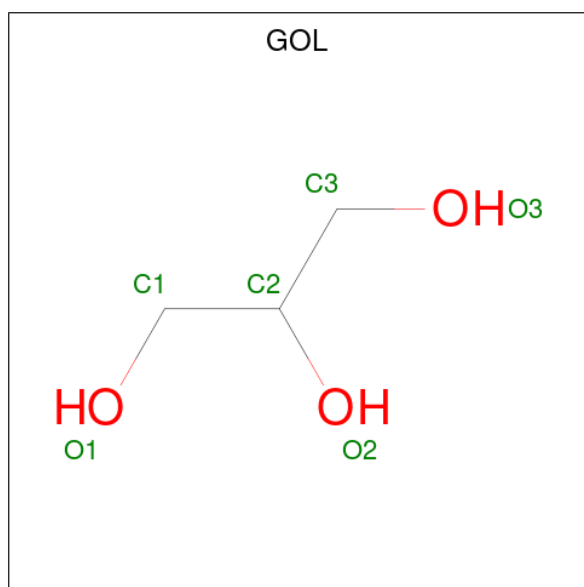
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 29	C 21	N 3	O 5	0	0
2	B	1	Total 29	C 21	N 3	O 5	0	0

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

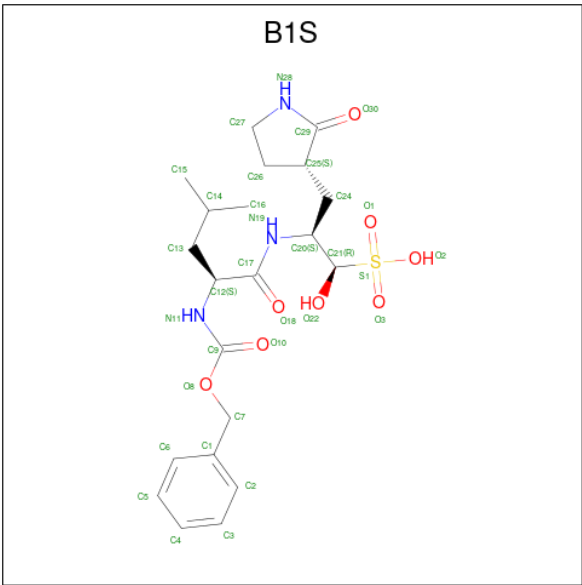
- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Cl	0	0
			3	3		
5	B	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		
6	B	1	Total	Mg	0	0
			1	1		

- Molecule 7 is (1R,2S)-2-({N-[(benzyloxy)carbonyl]-L-leucyl}amino)-1-hydroxy-3-[(3S)-2-oxo pyrrolidin-3-yl]propane-1-sulfonic acid (CCD ID: B1S) (formula: C₂₁H₃₁N₃O₈S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	N	O	0	0
			29	21	3	5		

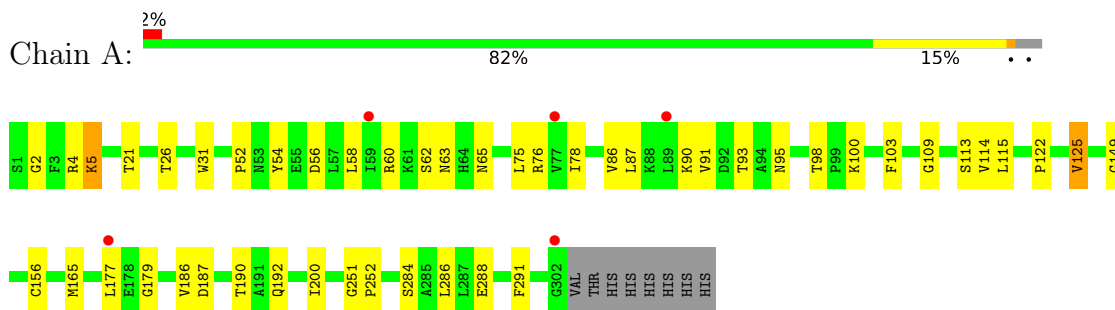
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	131	Total	O	0	7
			138	138		
8	B	121	Total	O	0	3
			124	124		
8	C	75	Total	O	0	3
			78	78		

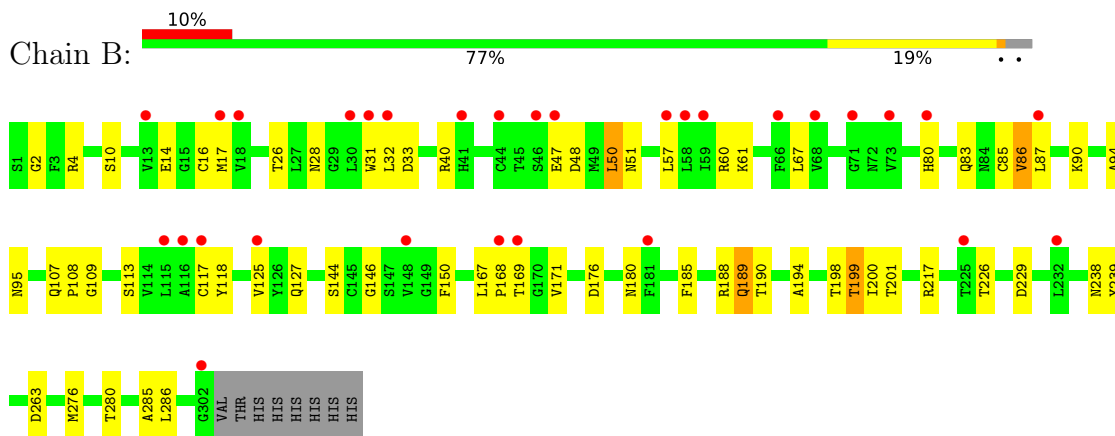
3 Residue-property plots [i](#)

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.

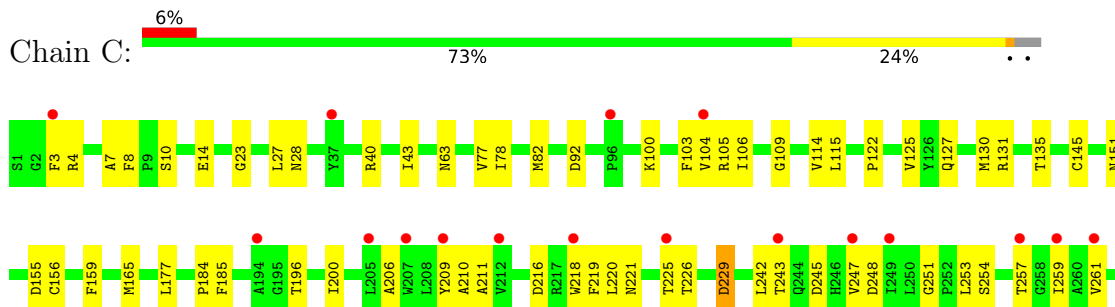
- Molecule 1: 3C-like proteinase

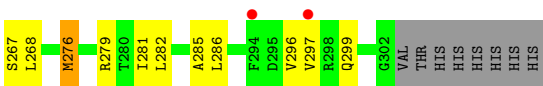


- Molecule 1: 3C-like proteinase



- Molecule 1: 3C-like proteinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.82Å 101.82Å 160.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.56 – 2.15 48.56 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.56-2.15) 99.9 (48.56-2.15)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.214 , 0.300 0.225 , 0.304	Depositor DCC
R_{free} test set	2733 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	0.798	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7430	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% 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: CL, B1S, PEG, GOL, MG, K36

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	1.04	1/2377 (0.0%)	1.38	1/3230 (0.0%)
1	B	1.06	1/2385 (0.0%)	1.40	5/3241 (0.2%)
1	C	1.05	0/2375	1.42	6/3227 (0.2%)
All	All	1.05	2/7137 (0.0%)	1.40	12/9698 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	VAL	C-O	5.62	1.29	1.24
1	B	176	ASP	C-O	5.26	1.29	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	GLY	CA-C-O	-6.31	118.12	122.22
1	B	201	THR	CA-CB-OG1	-6.25	100.22	109.60
1	B	199	THR	CA-CB-OG1	-5.63	101.16	109.60
1	B	2	GLY	CA-C-O	-5.41	118.71	122.22
1	C	23	GLY	CA-C-N	5.33	127.68	120.38
1	C	23	GLY	C-N-CA	5.33	127.68	120.38
1	B	263	ASP	CA-CB-CG	5.32	117.92	112.60
1	C	28	ASN	CA-C-N	5.25	125.80	122.18
1	C	28	ASN	C-N-CA	5.25	125.80	122.18
1	C	184	PRO	CA-C-O	-5.22	117.45	121.31
1	B	60	ARG	N-CA-C	-5.15	105.85	111.82
1	C	92	ASP	CB-CA-C	-5.00	102.49	110.79

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	2326	0	2273	26	0
1	B	2333	0	2280	51	0
1	C	2324	0	2266	57	0
2	A	29	0	29	0	0
2	B	29	0	29	0	0
3	A	7	0	10	0	0
4	A	6	0	8	0	0
5	A	3	0	0	0	0
5	B	1	0	0	1	0
5	C	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	C	29	0	0	0	0
8	A	138	0	0	0	0
8	B	124	0	0	2	0
8	C	78	0	0	0	0
All	All	7430	0	6895	127	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 (127) 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:C:257:THR:HG22	1:C:259:ILE:HG13	1.23	1.16
1:C:27:LEU:HD11	1:C:145:CYS:HB2	1.52	0.92
1:B:50:LEU:HD23	1:B:50:LEU:H	1.34	0.92
1:C:257:THR:CG2	1:C:259:ILE:HG13	2.08	0.81
1:B:28:ASN:O	1:B:146:GLY:HA3	1.81	0.81
1:A:103:PHE:CE1	1:A:177:LEU:HD23	2.22	0.73
1:B:31:TRP:CE2	1:B:95:ASN:HB2	2.26	0.70
1:C:104:VAL:HG12	1:C:105:ARG:H	1.57	0.70
1:B:31:TRP:CZ2	1:B:95:ASN:N	2.61	0.69
1:B:226:THR:OG1	1:B:229:ASP:HB2	1.91	0.69
1:C:218:TRP:CE3	1:C:279:ARG:HD2	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:VAL:HG12	1:C:105:ARG:N	2.08	0.69
1:B:31:TRP:CD2	1:B:95:ASN:HB2	2.29	0.68
1:B:198:THR:HG22	1:B:238:ASN:OD1	1.92	0.67
1:C:210:ALA:HB2	1:C:296:VAL:HG13	1.78	0.66
1:B:57:LEU:O	1:B:61:LYS:HG3	1.96	0.66
1:B:40:ARG:HD3	1:B:85:CYS:HA	1.79	0.65
1:B:31:TRP:CZ2	1:B:95:ASN:HA	2.33	0.64
1:B:31:TRP:CZ2	1:B:95:ASN:CA	2.81	0.64
1:B:50:LEU:H	1:B:50:LEU:CD2	1.99	0.63
1:B:50:LEU:HD23	1:B:50:LEU:N	2.12	0.62
1:C:100:LYS:O	1:C:156:CYS:HA	2.00	0.62
1:A:56:ASP:OD1	1:A:60:ARG:NH2	2.33	0.62
1:C:104:VAL:CG1	1:C:105:ARG:H	2.12	0.62
1:C:63:ASN:ND2	1:C:78:ILE:O	2.33	0.60
1:B:285:ALA:HB3	1:C:285:ALA:HB3	1.84	0.60
1:B:127:GLN:O	1:C:4:ARG:NH2	2.37	0.58
1:A:5:LYS:HD3	1:A:291:PHE:CZ	2.39	0.57
1:B:113:SER:O	1:B:150:PHE:CD1	2.57	0.57
1:B:48:ASP:HA	1:B:50:LEU:HD21	1.87	0.57
1:B:31:TRP:CH2	1:B:95:ASN:HA	2.39	0.57
1:B:47:GLU:O	1:B:50:LEU:HD21	2.05	0.56
1:B:169:THR:HG23	1:B:171:VAL:HG22	1.87	0.56
1:C:115:LEU:HD11	1:C:122:PRO:HB3	1.87	0.55
1:B:51:ASN:HA	1:B:188:ARG:HG2	1.87	0.55
1:A:86:VAL:HG13	1:A:179:GLY:HA2	1.88	0.55
1:A:103:PHE:CE1	1:A:177:LEU:CD2	2.90	0.55
1:B:276:MET:HE3	1:B:280:THR:HA	1.89	0.54
1:A:284:SER:OG	1:A:288:GLU:OE2	2.21	0.54
1:C:104:VAL:CG1	1:C:105:ARG:N	2.69	0.54
1:C:226:THR:OG1	1:C:229:ASP:OD2	2.24	0.53
1:C:209:TYR:HB3	1:C:253:LEU:HD22	1.90	0.53
1:C:131:ARG:HG2	1:C:135:THR:O	2.09	0.53
1:C:242:LEU:HG	1:C:247:VAL:HG22	1.91	0.53
1:B:167:LEU:HB3	1:B:168:PRO:HD2	1.90	0.53
1:C:100:LYS:HG2	1:C:156:CYS:HB2	1.90	0.53
1:C:165:MET:HE1	1:C:185:PHE:HB3	1.91	0.52
1:B:285:ALA:CB	1:C:285:ALA:HB3	2.40	0.52
1:A:76:ARG:O	1:A:91:VAL:HA	2.10	0.52
1:B:109:GLY:HA2	1:B:200:ILE:HD13	1.91	0.52
1:A:95:ASN:HB3	1:A:98:THR:OG1	2.10	0.51
1:B:217:ARG:NH2	8:B:505:HOH:O	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:103:PHE:CE1	1:C:177:LEU:HD23	2.46	0.50
1:B:10:SER:OG	1:B:14:GLU:OE2	2.29	0.50
1:B:169:THR:HG22	8:B:590:HOH:O	2.11	0.49
1:C:114:VAL:O	1:C:125:VAL:HA	2.11	0.49
1:B:40:ARG:CD	1:B:85:CYS:HA	2.42	0.49
1:C:8:PHE:HE2	1:C:151:ASN:OD1	1.96	0.49
1:C:10:SER:O	1:C:14:GLU:HG3	2.12	0.49
1:C:253:LEU:O	1:C:257:THR:OG1	2.23	0.49
1:C:211:ALA:O	1:C:216:ASP:HB2	2.12	0.49
1:C:40:ARG:HB2	1:C:82:MET:HE2	1.95	0.48
1:B:286:LEU:HD21	1:C:285:ALA:HB2	1.94	0.48
1:C:276:MET:CE	1:C:281:ILE:HD12	2.43	0.48
1:A:103:PHE:CZ	1:A:177:LEU:HD23	2.48	0.48
1:C:286:LEU:C	1:C:286:LEU:HD12	2.38	0.48
1:B:199:THR:HG21	1:B:239:TYR:CZ	2.49	0.48
1:C:155:ASP:OD1	1:C:155:ASP:N	2.39	0.48
1:A:109:GLY:HA2	1:A:200:ILE:HD13	1.95	0.47
1:C:109:GLY:HA2	1:C:200:ILE:HD13	1.95	0.47
1:C:251:GLY:O	1:C:254:SER:HB3	2.14	0.47
1:B:118:TYR:CE2	1:B:144:SER:HB3	2.50	0.47
1:C:159:PHE:HB3	1:C:177:LEU:HD22	1.96	0.47
1:C:243:THR:O	1:C:247:VAL:HG23	2.14	0.47
1:A:52:PRO:HG2	1:A:54:TYR:CE2	2.51	0.46
1:B:16:CYS:HA	1:B:95:ASN:ND2	2.31	0.46
1:B:16:CYS:HA	1:B:95:ASN:HD21	1.79	0.46
1:B:167:LEU:HD23	1:B:194:ALA:H	1.80	0.46
1:A:21:THR:HG23	1:A:26:THR:OG1	2.15	0.46
1:A:31:TRP:CE3	1:A:75:LEU:HD11	2.50	0.46
1:A:58:LEU:HD13	1:A:87:LEU:HD21	1.98	0.46
1:B:125:VAL:HG12	1:C:7:ALA:HB3	1.97	0.46
1:A:31:TRP:CZ2	1:A:95:ASN:HA	2.51	0.46
1:B:4:ARG:NH1	1:C:127:GLN:O	2.49	0.45
1:B:167:LEU:CB	1:B:168:PRO:HD2	2.46	0.45
1:C:3:PHE:CE1	1:C:299:GLN:NE2	2.85	0.45
1:B:80:HIS:CD2	1:B:80:HIS:O	2.69	0.45
1:B:169:THR:CG2	1:B:171:VAL:HG22	2.47	0.45
1:B:31:TRP:CE2	1:B:95:ASN:N	2.85	0.45
1:C:245:ASP:O	1:C:248:ASP:HB2	2.17	0.45
1:C:165:MET:CE	1:C:185:PHE:HB3	2.47	0.44
1:B:189:GLN:HE21	1:B:189:GLN:HB3	1.52	0.44
1:C:40:ARG:O	1:C:43:ILE:HG12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:243:THR:O	1:C:247:VAL:N	2.50	0.44
1:C:63:ASN:HB3	1:C:77:VAL:O	2.18	0.44
1:B:67:LEU:HG	5:B:402:CL:CL	2.55	0.44
1:A:100:LYS:O	1:A:156:CYS:HA	2.18	0.43
1:B:32:LEU:O	1:B:33:ASP:C	2.60	0.43
1:C:27:LEU:HD12	1:C:27:LEU:HA	1.67	0.43
1:A:63:ASN:ND2	1:A:78:ILE:C	2.77	0.43
1:A:115:LEU:HD11	1:A:122:PRO:HB3	2.00	0.43
1:B:17:MET:HG3	1:B:117:CYS:SG	2.59	0.43
1:A:62:SER:N	1:A:65:ASN:OD1	2.51	0.43
1:C:159:PHE:CB	1:C:177:LEU:HD22	2.49	0.43
1:B:31:TRP:CZ2	1:B:94:ALA:C	2.97	0.43
1:C:206:ALA:O	1:C:296:VAL:HG21	2.19	0.42
1:C:219:PHE:O	1:C:267:SER:OG	2.34	0.42
1:C:220:LEU:N	1:C:220:LEU:HD23	2.34	0.42
1:A:90:LYS:HE3	1:B:180:ASN:OD1	2.19	0.42
1:A:113:SER:O	1:A:149:GLY:HA2	2.20	0.42
1:C:210:ALA:HB2	1:C:296:VAL:CG1	2.48	0.42
1:C:221:ASN:HD22	1:C:221:ASN:HA	1.70	0.42
1:A:165:MET:HE1	1:A:192:GLN:NE2	2.35	0.42
1:C:8:PHE:CE2	1:C:151:ASN:OD1	2.73	0.42
1:A:114:VAL:O	1:A:125:VAL:HA	2.19	0.42
1:A:286:LEU:C	1:A:286:LEU:HD12	2.45	0.41
1:B:167:LEU:HD23	1:B:194:ALA:N	2.34	0.41
1:B:167:LEU:HD21	1:B:185:PHE:CE1	2.55	0.41
1:C:276:MET:CE	1:C:281:ILE:CD1	2.99	0.41
1:A:54:TYR:OH	1:A:187:ASP:O	2.38	0.41
1:C:209:TYR:OH	1:C:261:VAL:HA	2.20	0.41
1:B:86:VAL:O	1:B:86:VAL:HG12	2.20	0.40
1:A:251:GLY:N	1:A:252:PRO:HD2	2.37	0.40
1:B:107:GLN:O	1:B:108:PRO:C	2.63	0.40
1:C:106:ILE:HD11	1:C:130:MET:HB2	2.03	0.40
1:C:3:PHE:HB3	1:C:282:LEU:HD22	2.03	0.40
1:C:297:VAL:HG12	1:C:297:VAL:O	2.20	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	300/310 (97%)	284 (95%)	16 (5%)	0	100	100
1	B	300/310 (97%)	286 (95%)	14 (5%)	0	100	100
1	C	300/310 (97%)	283 (94%)	17 (6%)	0	100	100
All	All	900/930 (97%)	853 (95%)	47 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	258/267 (97%)	253 (98%)	5 (2%)	50	56
1	B	259/267 (97%)	251 (97%)	8 (3%)	35	37
1	C	257/267 (96%)	252 (98%)	5 (2%)	50	56
All	All	774/801 (97%)	756 (98%)	18 (2%)	44	49

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	5	LYS
1	A	93	THR
1	A	125	VAL
1	A	190	THR

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Mol	Chain	Res	Type
1	B	26	THR
1	B	50	LEU
1	B	83	GLN
1	B	86	VAL
1	B	87	LEU
1	B	90	LYS
1	B	189	GLN
1	B	190	THR
1	C	196	THR
1	C	225	THR
1	C	229	ASP
1	C	268	LEU
1	C	276	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
1	A	246	HIS
1	A	273	GLN
1	B	142	ASN
1	B	189	GLN
1	C	19	GLN
1	C	84	ASN
1	C	151	ASN
1	C	244	GLN
1	C	273	GLN
1	C	274	ASN

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 12 ligands modelled in this entry, 7 are monoatomic - leaving 5 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	K36	A	401	1	30,30,34	2.78	9 (30%)	35,39,47	1.88	6 (17%)
3	PEG	A	402	-	6,6,6	0.17	0	5,5,5	0.07	0
4	GOL	A	403	-	5,5,5	0.08	0	5,5,5	0.28	0
7	B1S	C	401	1	30,30,34	2.12	9 (30%)	35,39,47	2.05	10 (28%)
2	K36	B	401	1	30,30,34	2.60	9 (30%)	35,39,47	1.48	5 (14%)

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	K36	A	401	1	-	2/27/37/45	0/2/2/2
3	PEG	A	402	-	-	2/4/4/4	-
4	GOL	A	403	-	-	2/4/4/4	-
7	B1S	C	401	1	-	6/27/37/45	0/2/2/2
2	K36	B	401	1	-	2/27/37/45	0/2/2/2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	K36	C29-N28	7.78	1.41	1.33
2	B	401	K36	C29-N28	7.71	1.41	1.33
7	C	401	B1S	C29-N28	7.30	1.41	1.33
2	A	401	K36	O8-C7	5.91	1.58	1.45
2	B	401	K36	O8-C9	5.40	1.46	1.35
2	A	401	K36	C24-C20	5.04	1.66	1.53
2	B	401	K36	C9-N11	4.99	1.47	1.34
7	C	401	B1S	C7-C1	4.54	1.61	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	K36	C17-N19	4.51	1.43	1.34
2	B	401	K36	C17-N19	4.49	1.43	1.34
2	A	401	K36	C12-C17	4.17	1.63	1.52
2	A	401	K36	O8-C9	3.99	1.43	1.35
2	A	401	K36	C20-N19	-3.40	1.40	1.46
2	B	401	K36	C21-C20	3.35	1.58	1.52
2	B	401	K36	C20-N19	3.20	1.53	1.46
2	B	401	K36	O8-C7	3.09	1.52	1.45
7	C	401	B1S	C24-C25	-2.97	1.47	1.53
7	C	401	B1S	C25-C29	2.96	1.59	1.52
2	B	401	K36	C12-C17	-2.76	1.45	1.52
2	A	401	K36	C7-C1	2.49	1.56	1.50
2	A	401	K36	C6-C1	2.39	1.43	1.38
7	C	401	B1S	C4-C3	2.31	1.43	1.38
7	C	401	B1S	C5-C6	2.27	1.42	1.38
7	C	401	B1S	C13-C12	2.20	1.59	1.53
7	C	401	B1S	O8-C9	2.11	1.39	1.35
2	B	401	K36	C24-C20	2.05	1.58	1.53
7	C	401	B1S	O10-C9	2.04	1.25	1.21

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	K36	O8-C7-C1	5.11	121.80	109.40
2	A	401	K36	O8-C9-N11	4.97	121.07	110.45
7	C	401	B1S	C26-C25-C29	-4.85	96.21	102.87
7	C	401	B1S	O8-C7-C1	4.70	120.81	109.40
2	B	401	K36	O8-C9-N11	4.38	119.82	110.45
7	C	401	B1S	O30-C29-C25	-4.34	121.19	126.21
2	A	401	K36	C12-C17-N19	4.07	125.30	116.63
7	C	401	B1S	C24-C25-C29	3.90	121.25	112.81
7	C	401	B1S	C7-O8-C9	3.88	124.67	115.93
7	C	401	B1S	C27-N28-C29	-3.79	105.75	113.79
2	A	401	K36	O10-C9-N11	-3.67	118.84	124.86
2	B	401	K36	O8-C7-C1	3.51	117.92	109.40
2	B	401	K36	O8-C9-O10	-3.27	117.98	124.26
2	A	401	K36	O18-C17-N19	-3.04	117.51	122.96
7	C	401	B1S	C24-C20-C21	-2.32	106.49	111.30
2	B	401	K36	O18-C17-C12	-2.31	115.64	120.48
2	A	401	K36	O8-C9-O10	-2.27	119.91	124.26
7	C	401	B1S	O22-C21-C20	-2.24	106.10	111.73
7	C	401	B1S	C12-C17-N19	2.21	121.34	116.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	401	B1S	O8-C9-N11	2.09	114.93	110.45
2	B	401	K36	C17-C12-N11	-2.02	105.65	111.11

There are no chirality outliers.

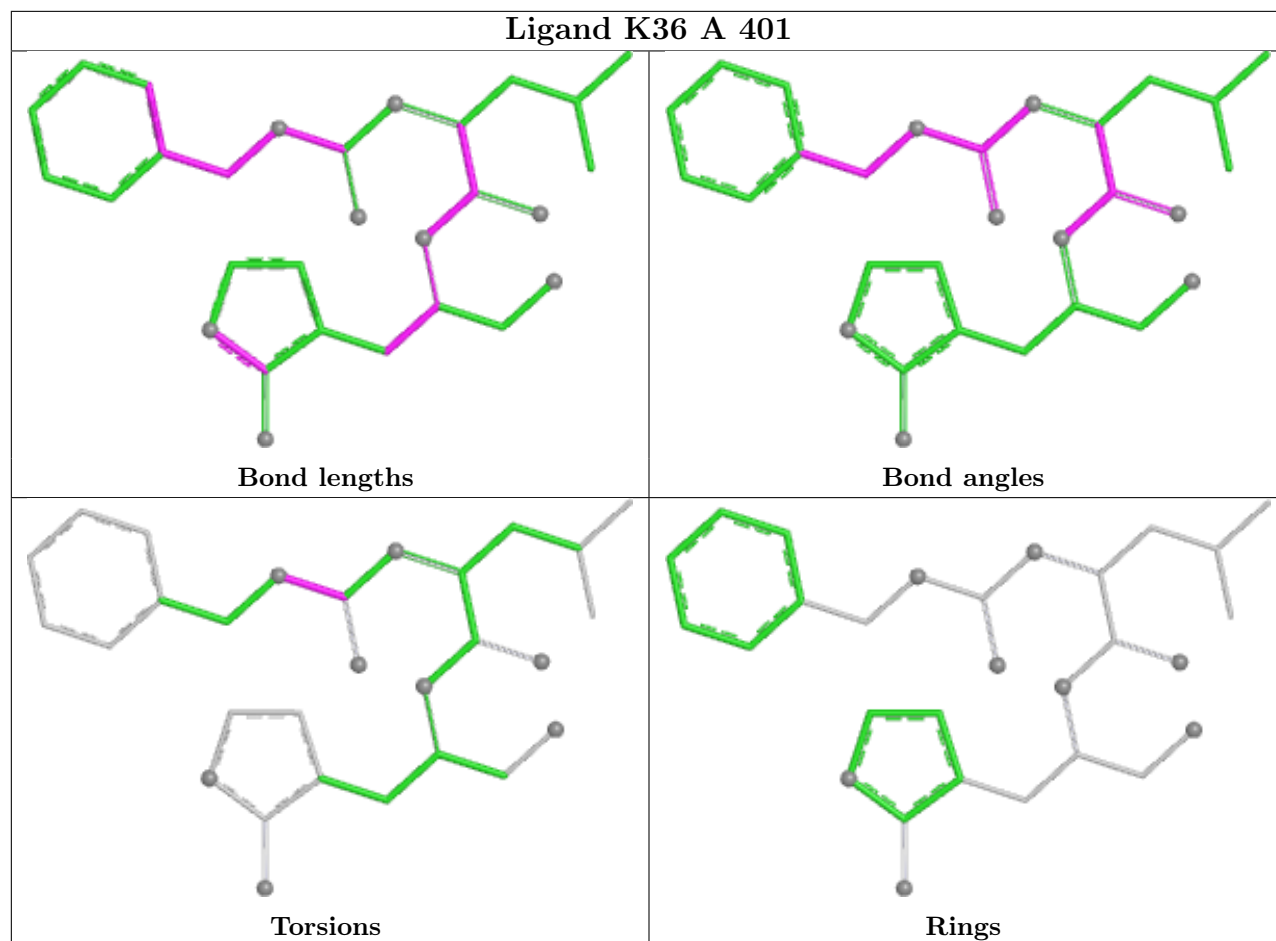
All (14) torsion outliers are listed below:

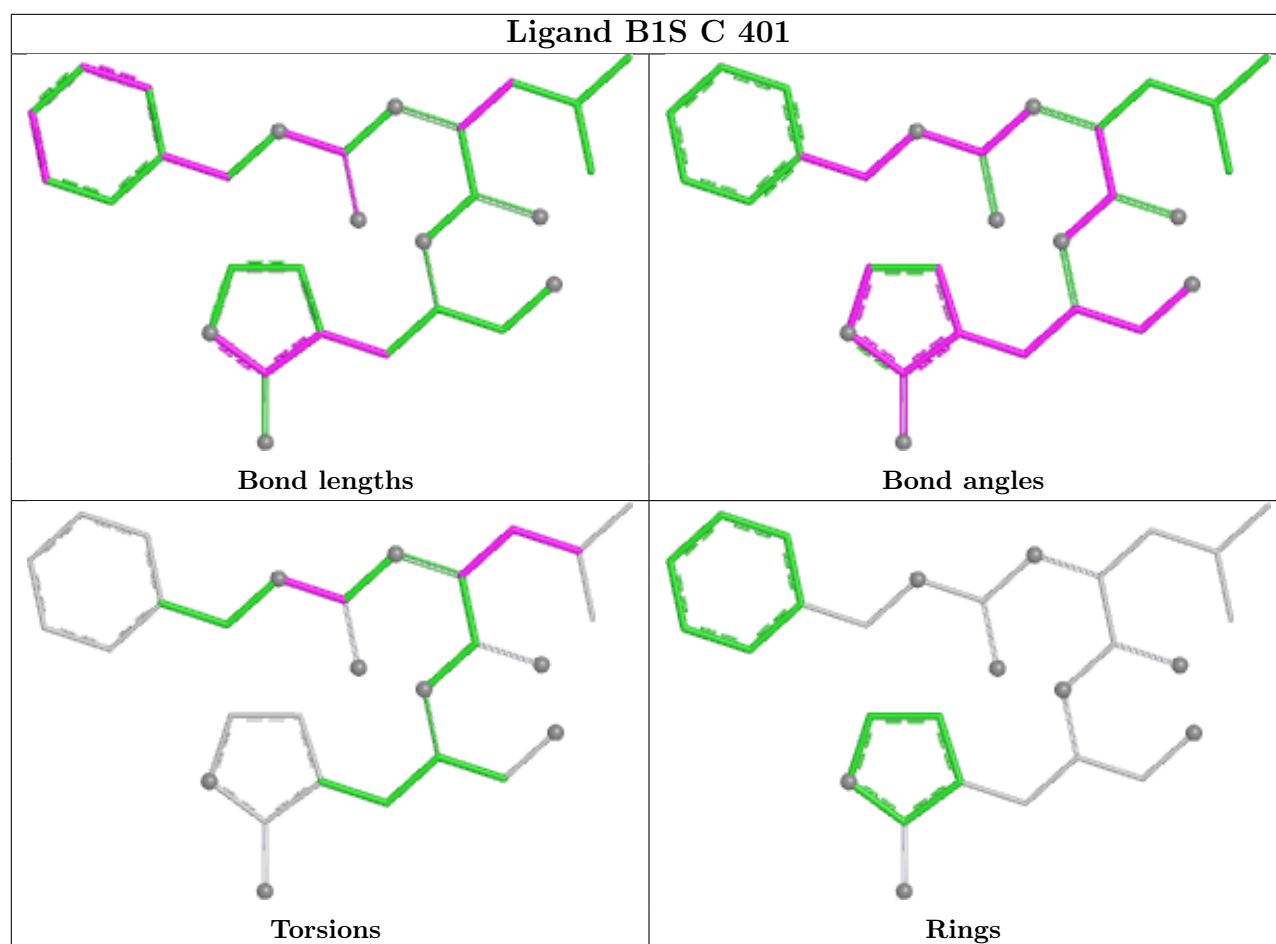
Mol	Chain	Res	Type	Atoms
2	A	401	K36	O10-C9-O8-C7
2	A	401	K36	N11-C9-O8-C7
2	B	401	K36	O10-C9-O8-C7
2	B	401	K36	N11-C9-O8-C7
4	A	403	GOL	O1-C1-C2-O2
7	C	401	B1S	C12-C13-C14-C16
7	C	401	B1S	N11-C9-O8-C7
4	A	403	GOL	O1-C1-C2-C3
3	A	402	PEG	O1-C1-C2-O2
7	C	401	B1S	O10-C9-O8-C7
7	C	401	B1S	C12-C13-C14-C15
3	A	402	PEG	O2-C3-C4-O4
7	C	401	B1S	N11-C12-C13-C14
7	C	401	B1S	C17-C12-C13-C14

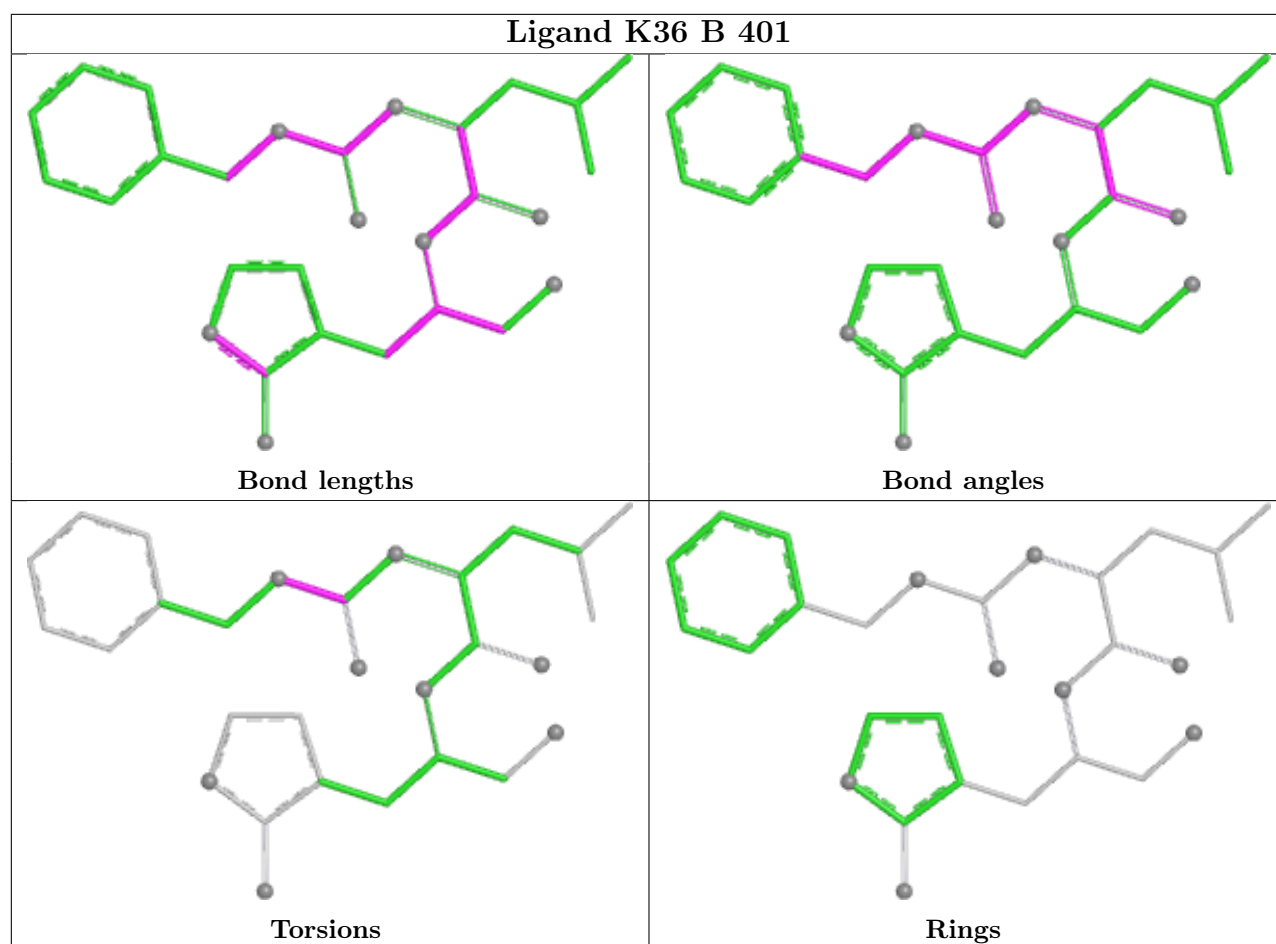
There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	302/310 (97%)	0.29	5 (1%) 69 73	41, 57, 95, 120	0
1	B	302/310 (97%)	0.75	30 (9%) 13 14	42, 64, 97, 123	0
1	C	302/310 (97%)	0.67	19 (6%) 26 29	45, 69, 99, 129	0
All	All	906/930 (97%)	0.57	54 (5%) 27 31	41, 63, 98, 129	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	148	VAL	4.3
1	B	116	ALA	3.6
1	B	73	VAL	3.5
1	B	302	GLY	3.4
1	B	168	PRO	3.2
1	C	209	TYR	3.1
1	B	68	VAL	3.0
1	C	297	VAL	2.8
1	B	46	SER	2.7
1	B	18	VAL	2.7
1	C	205	LEU	2.6
1	B	13	VAL	2.6
1	B	47	GLU	2.5
1	C	257	THR	2.5
1	B	41	HIS	2.5
1	B	71	GLY	2.5
1	C	207	TRP	2.5
1	C	194	ALA	2.4
1	C	3	PHE	2.4
1	C	243	THR	2.4
1	B	115	LEU	2.4
1	B	66	PHE	2.4
1	B	59	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	30	LEU	2.3
1	B	125	VAL	2.3
1	C	96	PRO	2.3
1	C	225	THR	2.3
1	C	104	VAL	2.3
1	B	232	LEU	2.3
1	A	59	ILE	2.2
1	B	169	THR	2.2
1	A	302	GLY	2.2
1	A	177	LEU	2.2
1	A	89	LEU	2.2
1	B	31	TRP	2.2
1	B	17	MET	2.2
1	C	259	ILE	2.2
1	A	77	VAL	2.2
1	C	261	VAL	2.2
1	B	117	CYS	2.2
1	B	80	HIS	2.1
1	B	44	CYS	2.1
1	B	225	THR	2.1
1	C	247	VAL	2.1
1	C	37	TYR	2.1
1	B	181	PHE	2.1
1	C	218	TRP	2.1
1	C	294	PHE	2.1
1	C	212	VAL	2.1
1	B	57	LEU	2.1
1	B	58	LEU	2.1
1	C	249	ILE	2.1
1	B	87	LEU	2.0
1	B	32	LEU	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

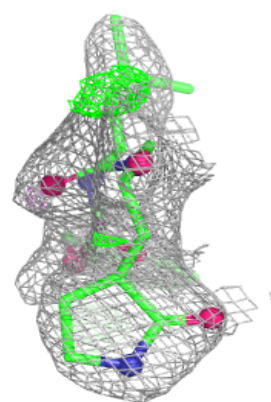
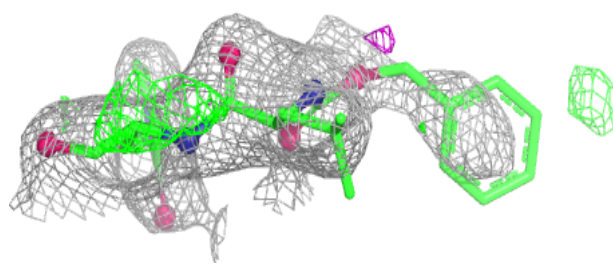
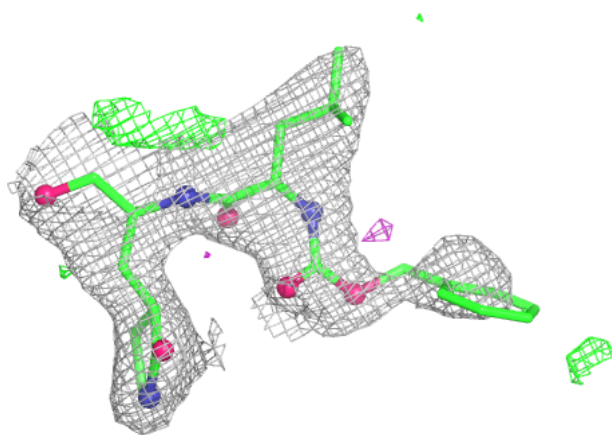
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PEG	A	402	7/7	0.75	0.13	70,82,85,87	0
5	CL	B	402	1/1	0.77	0.15	93,93,93,93	0
5	CL	A	405	1/1	0.79	0.18	90,90,90,90	0
6	MG	A	407	1/1	0.84	0.18	86,86,86,86	0
2	K36	B	401	29/33	0.85	0.13	51,67,112,116	0
4	GOL	A	403	6/6	0.87	0.12	72,77,80,82	0
5	CL	C	402	1/1	0.90	0.12	89,89,89,89	0
5	CL	A	404	1/1	0.90	0.14	97,97,97,97	0
2	K36	A	401	29/33	0.91	0.10	49,57,75,80	0
5	CL	A	406	1/1	0.92	0.08	87,87,87,87	0
7	B1S	C	401	29/33	0.95	0.08	45,50,68,68	0
6	MG	B	403	1/1	0.99	0.03	55,55,55,55	0

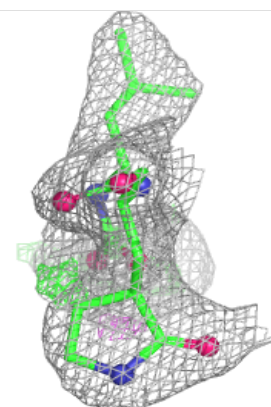
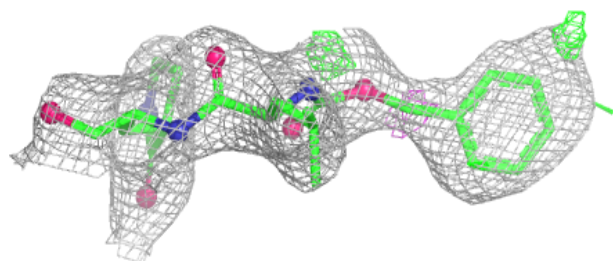
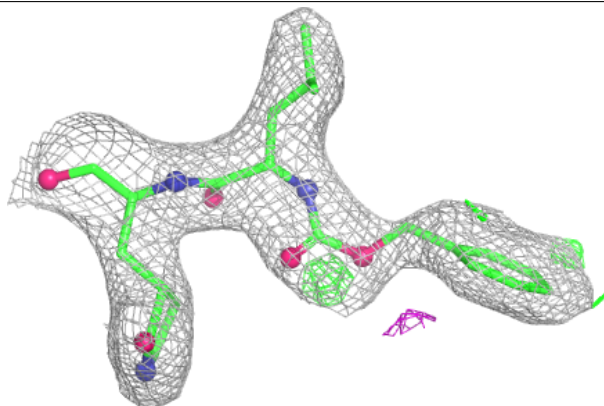
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around K36 B 401:

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

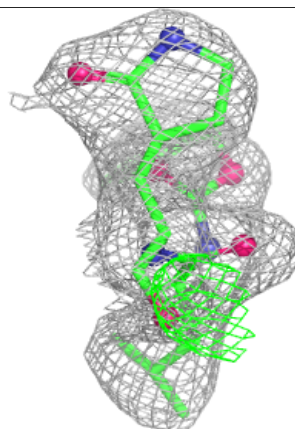
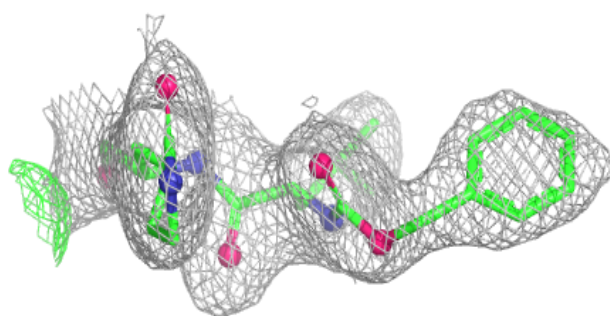
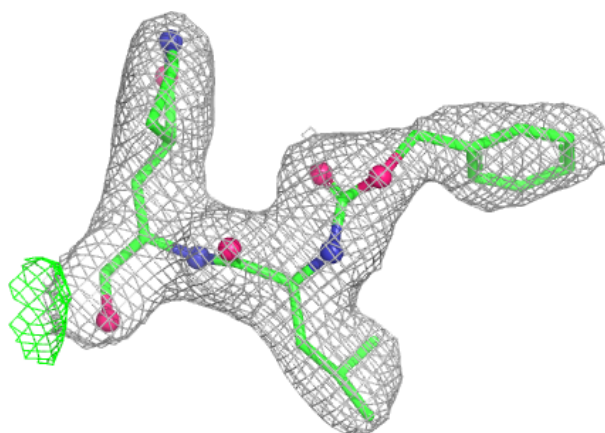
**Electron density around K36 A 401:**

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



Electron density around B1S C 401:

$2mF_o - DF_c$ (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.