



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 30, 2026 – 02:16 PM UTC

PDB ID : 4INU / pdb_00004inu
Title : Yeast 20S proteasome in complex with the vinyl sulfone LU112
Authors : Geurink, P.P.; van der Linden, W.A.; Mirabella, A.C.; Gallastegui, N.; de Bruin, G.; Blom, A.E.M.; Voges, M.J.; Mock, E.D.; Florea, B.I.; van der Marel, G.A.; Driessen, C.; van der Stelt, M.; Groll, M.; Overkleeft, H.S.; Kisselev, A.F.
Deposited on : 2013-01-06
Resolution : 3.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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](#) ①) 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

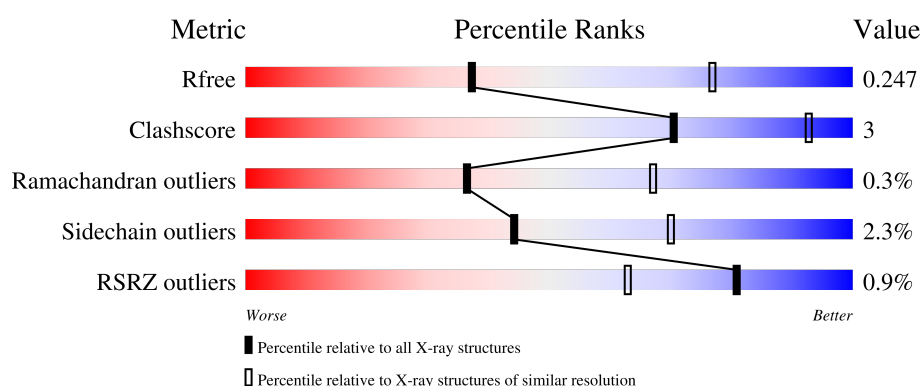
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	250	91% 8%
1	O	250	92% 8%
2	B	258	% 81% 13% • 5%
2	P	258	2% 82% 11% • 5%
3	C	254	86% 8% • 5%

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Mol	Chain	Length	Quality of chain
3	Q	254	 2% 87% 7% • 5%
4	D	260	 % 81% 12% • 7%
4	R	260	 % 80% 12% • 7%
5	E	234	 88% 11%
5	S	234	 3% 87% 12%
6	F	288	 % 76% 7% • 15%
6	T	288	 % 75% 9% • 15%
7	G	252	 87% 9% •
7	U	252	 89% 7% •
8	H	232	 89% 6% •
8	V	232	 89% 6% •
9	I	205	 86% 13%
9	W	205	 87% 13%
10	J	198	 2% 88% 11% •
10	X	198	 2% 90% 10% •
11	K	212	 3% 84% 15% •
11	Y	212	 2% 87% 12% •
12	L	222	 90% 9%
12	Z	222	 88% 12%
13	M	233	 90% 9%
13	a	233	 % 91% 9%
14	N	196	 93% 7% •
14	b	196	 95% 5% •

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 51078 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 Proteasome component Y7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			
1	O	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			

- Molecule 2 is a protein called Proteasome component Y13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			
2	P	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			

- Molecule 3 is a protein called Proteasome component PRE6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	241	Total	C	N	O	S	0	0	0
			1890	1181	331	374	4			
3	Q	241	Total	C	N	O	S	0	0	0
			1890	1181	331	374	4			

- Molecule 4 is a protein called Proteasome component PUP2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	242	Total	C	N	O	S	0	0	0
			1861	1162	314	378	7			
4	R	242	Total	C	N	O	S	0	0	0
			1861	1162	314	378	7			

- Molecule 5 is a protein called Proteasome component PRE5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	233	Total	C	N	O	S	0	0	0
			1795	1129	312	350	4			
5	S	233	Total	C	N	O	S	0	0	0
			1795	1129	312	350	4			

- Molecule 6 is a protein called Proteasome component C1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	244	Total	C	N	O	S	0	0	0
			1896	1205	330	357	4			
6	T	244	Total	C	N	O	S	0	0	0
			1896	1205	330	357	4			

- Molecule 7 is a protein called Proteasome component C7-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	243	Total	C	N	O	S	0	0	0
			1921	1221	322	370	8			
7	U	243	Total	C	N	O	S	0	0	0
			1921	1221	322	370	8			

- Molecule 8 is a protein called Proteasome component PUP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			
8	V	222	Total	C	N	O	S	0	0	0
			1684	1061	293	323	7			

- Molecule 9 is a protein called Proteasome component PUP3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	204	Total	C	N	O	S	0	0	0
			1581	1010	258	305	8			
9	W	204	Total	C	N	O	S	0	0	0
			1581	1010	258	305	8			

- Molecule 10 is a protein called Proteasome component C11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	198	Total	C	N	O	S	0	0	0
			1585	1005	269	305	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	X	198	Total	C	N	O	S	0	0	0
			1585	1005	269	305	6			

- Molecule 11 is a protein called Proteasome component PRE2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			

- Molecule 12 is a protein called Proteasome component C5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			
12	Z	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			

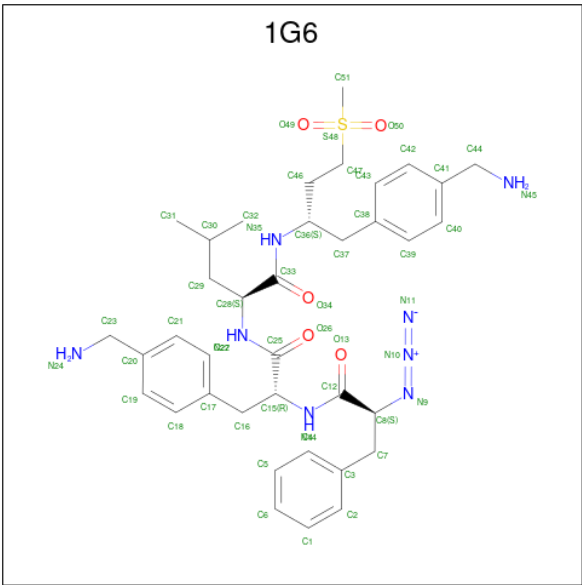
- Molecule 13 is a protein called Proteasome component PRE4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			
13	a	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			

- Molecule 14 is a protein called Proteasome component PRE3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			
14	b	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			

- Molecule 15 is N3Phe-Phe(4-NH₂CH₂)-Leu-Phe(4-NH₂CH₂)-methyl vinyl sulfone, bound form (CCD ID: 1G6) (formula: C₃₇H₅₀N₈O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
15	H	1	Total	C	N	O	S	0	0
			51	37	8	5	1		
15	K	1	Total	C	N	O	S	0	0
			51	37	8	5	1		
15	V	1	Total	C	N	O	S	0	0
			51	37	8	5	1		
15	Y	1	Total	C	N	O	S	0	0
			51	37	8	5	1		

- Molecule 16 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	60	Total	O	0	0
			60	60		
16	B	38	Total	O	0	0
			38	38		
16	C	42	Total	O	0	0
			42	42		
16	D	38	Total	O	0	0
			38	38		
16	E	20	Total	O	0	0
			20	20		
16	F	47	Total	O	0	0
			47	47		
16	G	61	Total	O	0	0
			61	61		
16	H	54	Total	O	0	0
			54	54		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	I	66	Total 66	O 66	0	0
16	J	50	Total 50	O 50	0	0
16	K	49	Total 49	O 49	0	0
16	L	56	Total 56	O 56	0	0
16	M	79	Total 79	O 79	0	0
16	N	54	Total 54	O 54	0	0
16	O	34	Total 34	O 34	0	0
16	P	29	Total 29	O 29	0	0
16	Q	29	Total 29	O 29	0	0
16	R	29	Total 29	O 29	0	0
16	S	18	Total 18	O 18	0	0
16	T	43	Total 43	O 43	0	0
16	U	59	Total 59	O 59	0	0
16	V	50	Total 50	O 50	0	0
16	W	54	Total 54	O 54	0	0
16	X	44	Total 44	O 44	0	0
16	Y	50	Total 50	O 50	0	0
16	Z	47	Total 47	O 47	0	0
16	a	74	Total 74	O 74	0	0
16	b	62	Total 62	O 62	0	0

3 Residue-property plots [i](#)

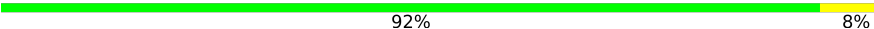
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Proteasome component Y7

Chain A: 



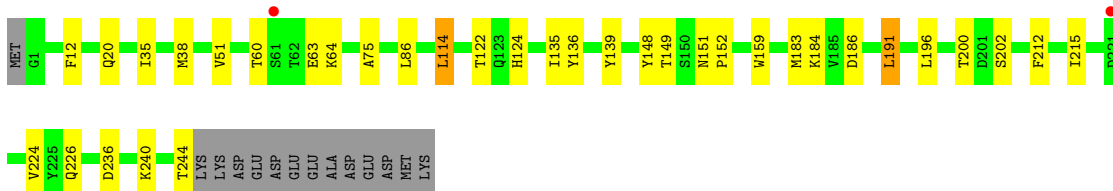
- Molecule 1: Proteasome component Y7

Chain O: 



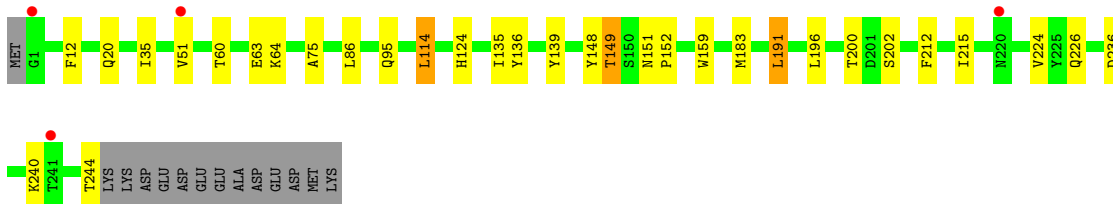
- Molecule 2: Proteasome component Y13

Chain B: 



- Molecule 2: Proteasome component Y13

Chain P: 



- Molecule 3: Proteasome component PRE6

Chain C:  86% 8% • 5%



- Molecule 3: Proteasome component PRE6

Chain Q:  87% 7% • 5%



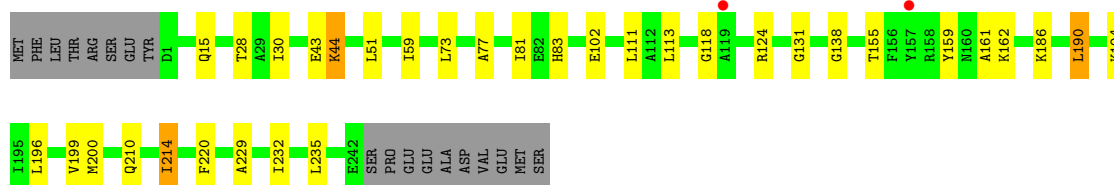
- Molecule 4: Proteasome component PUP2

Chain D:  81% 12% • 7%



- Molecule 4: Proteasome component PUP2

Chain R:  80% 12% • 7%



- Molecule 5: Proteasome component PRE5

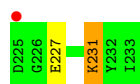
Chain E:  88% 11%



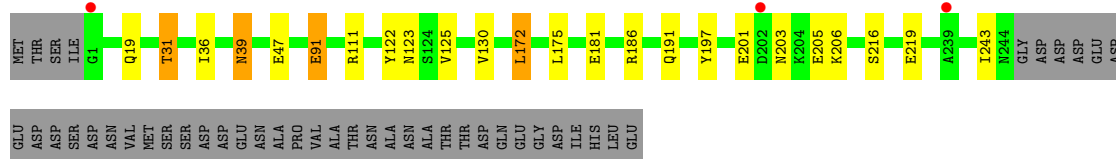
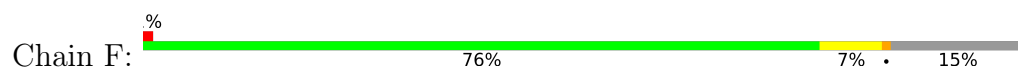
- Molecule 5: Proteasome component PRE5

Chain S:  87% 12% • 3%

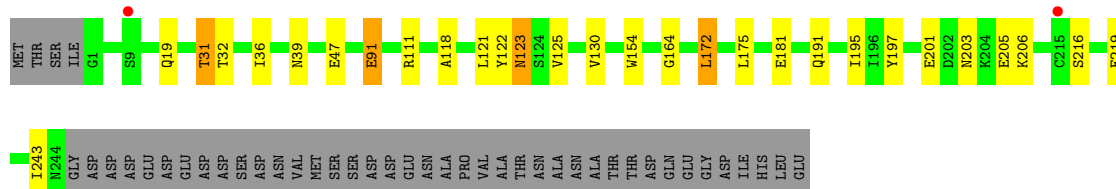
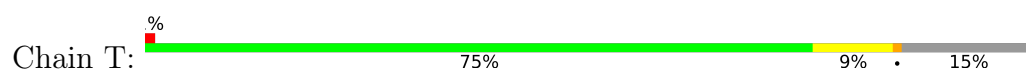




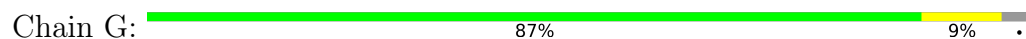
• Molecule 6: Proteasome component C1



• Molecule 6: Proteasome component C1



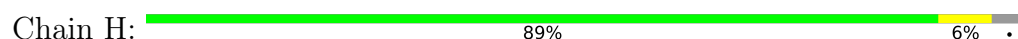
• Molecule 7: Proteasome component C7-alpha



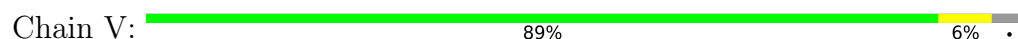
• Molecule 7: Proteasome component C7-alpha



• Molecule 8: Proteasome component PUP1

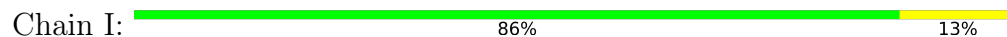


• Molecule 8: Proteasome component PUP1

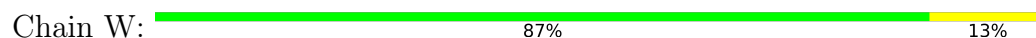




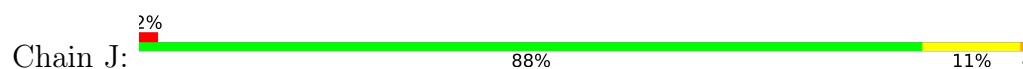
- Molecule 9: Proteasome component PUP3



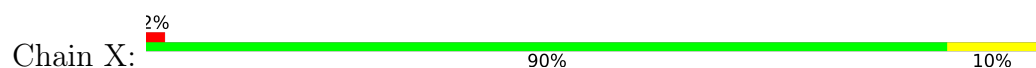
- Molecule 9: Proteasome component PUP3



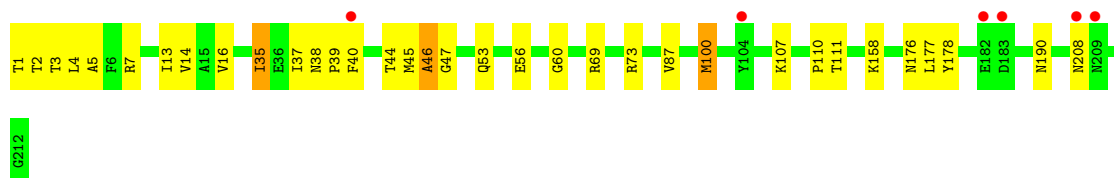
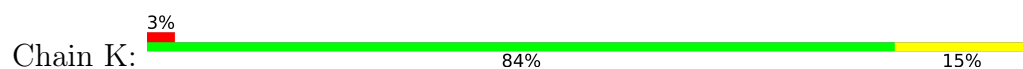
- Molecule 10: Proteasome component C11



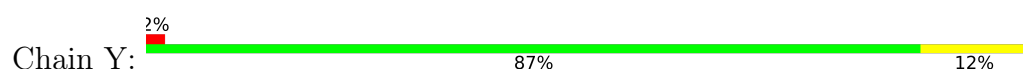
- Molecule 10: Proteasome component C11



- Molecule 11: Proteasome component PRE2



- Molecule 11: Proteasome component PRE2



- Molecule 12: Proteasome component C5

Chain L:  90% 9%



- Molecule 12: Proteasome component C5

Chain Z:  88% 12%

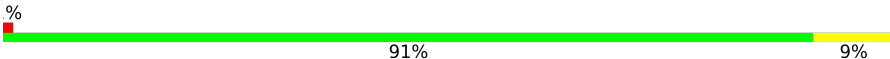


- Molecule 13: Proteasome component PRE4

Chain M:  90% 9%



- Molecule 13: Proteasome component PRE4

Chain a:  91% 9%



- Molecule 14: Proteasome component PRE3

Chain N:  93% 7%



- Molecule 14: Proteasome component PRE3

Chain b:  95% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	134.91Å 300.90Å 143.62Å 90.00° 112.53° 90.00°	Depositor
Resolution (Å)	15.00 – 3.10 15.00 – 3.10	Depositor EDS
% Data completeness (in resolution range)	92.3 (15.00-3.10) 91.7 (15.00-3.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.08 (at 3.12Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.238 0.226 , 0.247	Depositor DCC
R_{free} test set	8746 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	75.3	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 59.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	51078	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% 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: 1G6

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	0.49	0/1952	0.76	0/2642
1	O	0.49	0/1952	0.77	0/2642
2	B	0.47	0/1934	0.74	0/2618
2	P	0.47	0/1934	0.73	0/2618
3	C	0.47	0/1919	0.76	0/2598
3	Q	0.48	0/1919	0.77	0/2598
4	D	0.53	0/1886	0.72	0/2541
4	R	0.53	0/1886	0.72	0/2541
5	E	0.48	0/1823	0.73	2/2463 (0.1%)
5	S	0.47	0/1823	0.73	2/2463 (0.1%)
6	F	0.59	0/1936	0.78	0/2614
6	T	0.59	0/1936	0.77	0/2614
7	G	0.49	0/1959	0.74	0/2652
7	U	0.49	0/1959	0.74	0/2652
8	H	0.62	0/1715	0.73	0/2326
8	V	0.62	0/1715	0.74	0/2326
9	I	0.45	0/1611	0.73	0/2174
9	W	0.46	0/1611	0.72	0/2174
10	J	0.48	0/1613	0.73	0/2173
10	X	0.49	0/1613	0.74	0/2173
11	K	0.59	0/1681	0.74	0/2274
11	Y	0.59	0/1681	0.75	0/2274
12	L	0.51	0/1795	0.73	0/2420
12	Z	0.51	0/1795	0.72	0/2420
13	M	0.47	0/1855	0.73	0/2514
13	a	0.47	0/1855	0.73	0/2514
14	N	0.52	0/1541	0.73	0/2087
14	b	0.52	0/1541	0.74	0/2087
All	All	0.51	0/50440	0.74	4/68192 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	S	147	GLN	CA-C-N	5.42	125.75	119.47
5	S	147	GLN	C-N-CA	5.42	125.75	119.47
5	E	147	GLN	CA-C-N	5.16	125.45	119.47
5	E	147	GLN	C-N-CA	5.16	125.45	119.47

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	1915	0	1929	11	0
1	O	1915	0	1929	9	0
2	B	1904	0	1904	18	0
2	P	1904	0	1904	17	0
3	C	1890	0	1903	15	0
3	Q	1890	0	1903	11	0
4	D	1861	0	1839	17	0
4	R	1861	0	1839	19	0
5	E	1795	0	1800	14	0
5	S	1795	0	1800	15	0
6	F	1896	0	1889	14	0
6	T	1896	0	1889	17	0
7	G	1921	0	1913	10	0
7	U	1921	0	1913	9	0
8	H	1684	0	1687	7	0
8	V	1684	0	1687	7	0
9	I	1581	0	1574	16	0
9	W	1581	0	1574	16	0
10	J	1585	0	1590	14	0
10	X	1585	0	1590	12	0
11	K	1644	0	1594	17	0
11	Y	1644	0	1594	21	0
12	L	1757	0	1711	11	0
12	Z	1757	0	1711	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	1824	0	1832	14	0
13	a	1824	0	1832	12	0
14	N	1512	0	1481	8	0
14	b	1512	0	1481	6	0
15	H	51	0	49	0	0
15	K	51	0	49	1	0
15	V	51	0	49	2	0
15	Y	51	0	49	8	0
16	A	60	0	0	0	0
16	B	38	0	0	0	0
16	C	42	0	0	0	0
16	D	38	0	0	0	0
16	E	20	0	0	0	0
16	F	47	0	0	0	0
16	G	61	0	0	0	0
16	H	54	0	0	0	0
16	I	66	0	0	0	0
16	J	50	0	0	1	0
16	K	49	0	0	0	0
16	L	56	0	0	0	0
16	M	79	0	0	0	0
16	N	54	0	0	0	0
16	O	34	0	0	0	0
16	P	29	0	0	0	0
16	Q	29	0	0	0	0
16	R	29	0	0	0	0
16	S	18	0	0	0	0
16	T	43	0	0	0	0
16	U	59	0	0	0	0
16	V	50	0	0	0	0
16	W	54	0	0	0	0
16	X	44	0	0	0	0
16	Y	50	0	0	0	0
16	Z	47	0	0	0	0
16	a	74	0	0	0	0
16	b	62	0	0	0	0
All	All	51078	0	49488	333	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 3.

The worst 5 of 333 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)
11:Y:1:THR:N	15:Y:301:1G6:H5	1.59	1.17
15:Y:301:1G6:H3	15:Y:301:1G6:H22	1.52	0.89
11:Y:1:THR:CA	15:Y:301:1G6:H5	2.04	0.87
11:Y:1:THR:H1	15:Y:301:1G6:H5	1.31	0.85
11:Y:1:THR:N	15:Y:301:1G6:C47	2.40	0.85

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	248/250 (99%)	238 (96%)	9 (4%)	1 (0%)	30	61
1	O	248/250 (99%)	239 (96%)	8 (3%)	1 (0%)	30	61
2	B	242/258 (94%)	236 (98%)	5 (2%)	1 (0%)	30	61
2	P	242/258 (94%)	236 (98%)	5 (2%)	1 (0%)	30	61
3	C	239/254 (94%)	232 (97%)	5 (2%)	2 (1%)	16	47
3	Q	239/254 (94%)	232 (97%)	5 (2%)	2 (1%)	16	47
4	D	240/260 (92%)	233 (97%)	4 (2%)	3 (1%)	9	35
4	R	240/260 (92%)	233 (97%)	6 (2%)	1 (0%)	30	61
5	E	231/234 (99%)	223 (96%)	8 (4%)	0	100	100
5	S	231/234 (99%)	223 (96%)	8 (4%)	0	100	100
6	F	242/288 (84%)	232 (96%)	10 (4%)	0	100	100
6	T	242/288 (84%)	232 (96%)	10 (4%)	0	100	100
7	G	241/252 (96%)	235 (98%)	6 (2%)	0	100	100
7	U	241/252 (96%)	236 (98%)	5 (2%)	0	100	100
8	H	220/232 (95%)	214 (97%)	6 (3%)	0	100	100
8	V	220/232 (95%)	213 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	I	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
9	W	202/205 (98%)	194 (96%)	8 (4%)	0	100	100
10	J	196/198 (99%)	189 (96%)	6 (3%)	1 (0%)	24	57
10	X	196/198 (99%)	188 (96%)	7 (4%)	1 (0%)	24	57
11	K	210/212 (99%)	200 (95%)	8 (4%)	2 (1%)	12	41
11	Y	210/212 (99%)	201 (96%)	8 (4%)	1 (0%)	24	57
12	L	220/222 (99%)	212 (96%)	8 (4%)	0	100	100
12	Z	220/222 (99%)	215 (98%)	5 (2%)	0	100	100
13	M	231/233 (99%)	220 (95%)	10 (4%)	1 (0%)	30	61
13	a	231/233 (99%)	217 (94%)	13 (6%)	1 (0%)	30	61
14	N	194/196 (99%)	186 (96%)	8 (4%)	0	100	100
14	b	194/196 (99%)	185 (95%)	9 (5%)	0	100	100
All	All	6312/6588 (96%)	6088 (96%)	205 (3%)	19 (0%)	36	67

5 of 19 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
11	K	46	ALA
11	K	39	PRO
3	Q	52	LEU
11	Y	39	PRO
1	A	2	THR

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	209/209 (100%)	206 (99%)	3 (1%)	59	76
1	O	209/209 (100%)	206 (99%)	3 (1%)	59	76
2	B	203/216 (94%)	197 (97%)	6 (3%)	36	65
2	P	203/216 (94%)	197 (97%)	6 (3%)	36	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	213/226 (94%)	209 (98%)	4 (2%)	50	73
3	Q	213/226 (94%)	208 (98%)	5 (2%)	44	70
4	D	198/215 (92%)	193 (98%)	5 (2%)	42	69
4	R	198/215 (92%)	193 (98%)	5 (2%)	42	69
5	E	192/193 (100%)	186 (97%)	6 (3%)	35	64
5	S	192/193 (100%)	186 (97%)	6 (3%)	35	64
6	F	201/239 (84%)	191 (95%)	10 (5%)	22	53
6	T	201/239 (84%)	192 (96%)	9 (4%)	24	56
7	G	207/210 (99%)	202 (98%)	5 (2%)	43	69
7	U	207/210 (99%)	203 (98%)	4 (2%)	50	73
8	H	181/190 (95%)	177 (98%)	4 (2%)	45	71
8	V	181/190 (95%)	177 (98%)	4 (2%)	45	71
9	I	172/173 (99%)	170 (99%)	2 (1%)	63	78
9	W	172/173 (99%)	170 (99%)	2 (1%)	63	78
10	J	175/175 (100%)	174 (99%)	1 (1%)	78	83
10	X	175/175 (100%)	173 (99%)	2 (1%)	65	78
11	K	169/169 (100%)	162 (96%)	7 (4%)	27	59
11	Y	169/169 (100%)	163 (96%)	6 (4%)	31	62
12	L	185/185 (100%)	182 (98%)	3 (2%)	55	75
12	Z	185/185 (100%)	183 (99%)	2 (1%)	65	78
13	M	199/199 (100%)	195 (98%)	4 (2%)	48	72
13	a	199/199 (100%)	193 (97%)	6 (3%)	36	65
14	N	162/162 (100%)	161 (99%)	1 (1%)	78	83
14	b	162/162 (100%)	161 (99%)	1 (1%)	78	83
All	All	5332/5522 (97%)	5210 (98%)	122 (2%)	44	70

5 of 122 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	M	70	LEU
11	Y	87	VAL
3	Q	147	GLN
11	Y	35	ILE
13	a	159	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 161 such sidechains are listed below:

Mol	Chain	Res	Type
5	S	116	GLN
11	Y	208	ASN
5	S	184	ASN
7	U	121	GLN
12	Z	165	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	$\# Z > 2$
15	1G6	Y	301	11	53,53,53	1.27	3 (5%)	66,71,71	1.79	15 (22%)
15	1G6	K	301	11	53,53,53	1.28	3 (5%)	66,71,71	1.59	10 (15%)
15	1G6	H	301	8	53,53,53	1.32	4 (7%)	66,71,71	1.47	8 (12%)
15	1G6	V	301	8	53,53,53	1.31	3 (5%)	66,71,71	1.71	10 (15%)

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
15	1G6	Y	301	11	-	13/53/53/53	0/3/3/3
15	1G6	K	301	11	-	13/53/53/53	0/3/3/3
15	1G6	H	301	8	-	17/53/53/53	0/3/3/3
15	1G6	V	301	8	-	19/53/53/53	0/3/3/3

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	V	301	1G6	C46-C47	4.66	1.57	1.52
15	H	301	1G6	C46-C47	4.54	1.57	1.52
15	K	301	1G6	C46-C47	4.20	1.56	1.52
15	Y	301	1G6	C47-S48	3.84	1.83	1.78
15	K	301	1G6	N11-N10	-3.37	1.16	1.23

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
15	V	301	1G6	C8-N9-N10	7.67	122.93	115.36
15	Y	301	1G6	C8-N9-N10	6.46	121.74	115.36
15	V	301	1G6	C15-C25-N27	6.33	130.13	116.63
15	K	301	1G6	C15-C25-N27	6.16	129.76	116.63
15	H	301	1G6	C15-C25-N27	5.94	129.30	116.63

There are no chirality outliers.

5 of 62 torsion outliers are listed below:

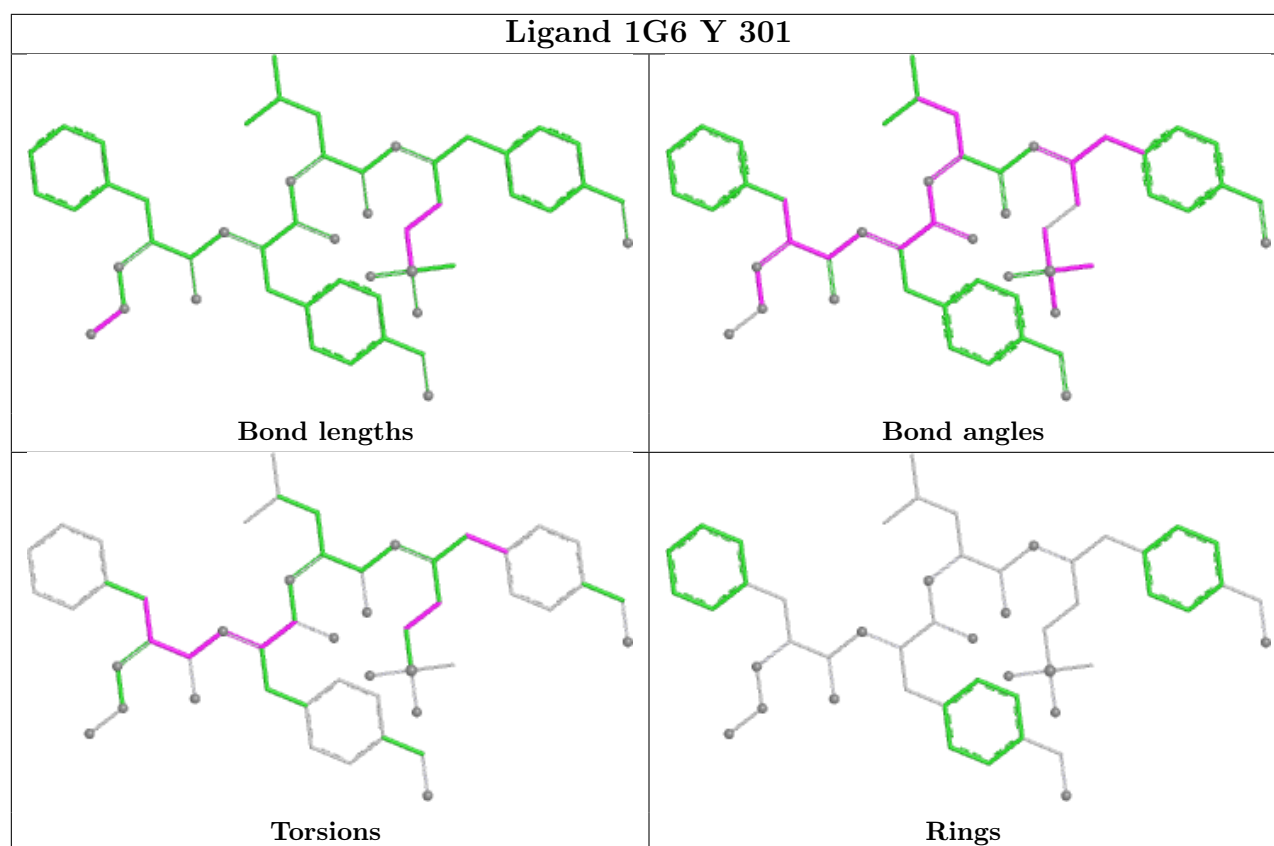
Mol	Chain	Res	Type	Atoms
15	H	301	1G6	O13-C12-N14-C15
15	H	301	1G6	C8-C12-N14-C15
15	H	301	1G6	C12-C8-N9-N10
15	H	301	1G6	C7-C8-N9-N10
15	H	301	1G6	C25-C15-N14-C12

There are no ring outliers.

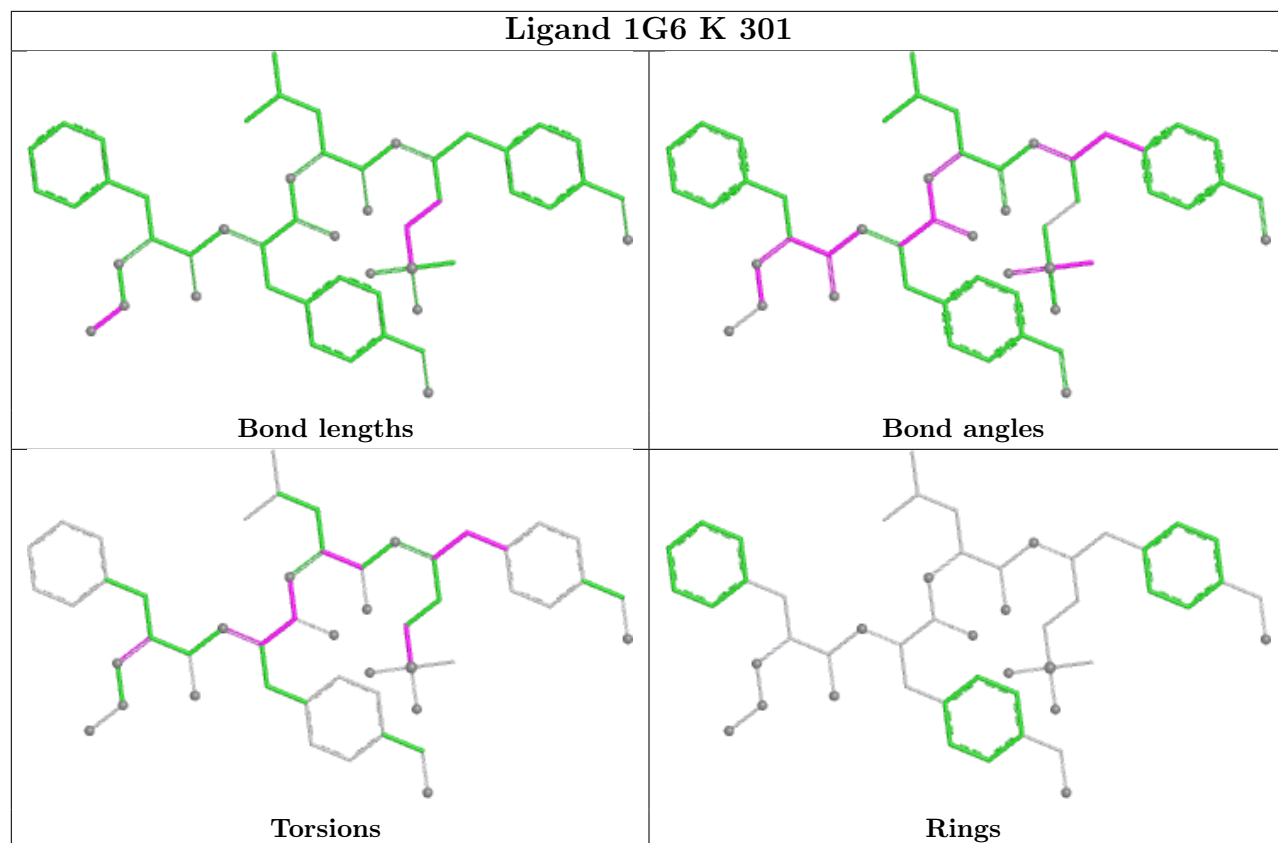
3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
15	Y	301	1G6	8	0
15	K	301	1G6	1	0
15	V	301	1G6	2	0

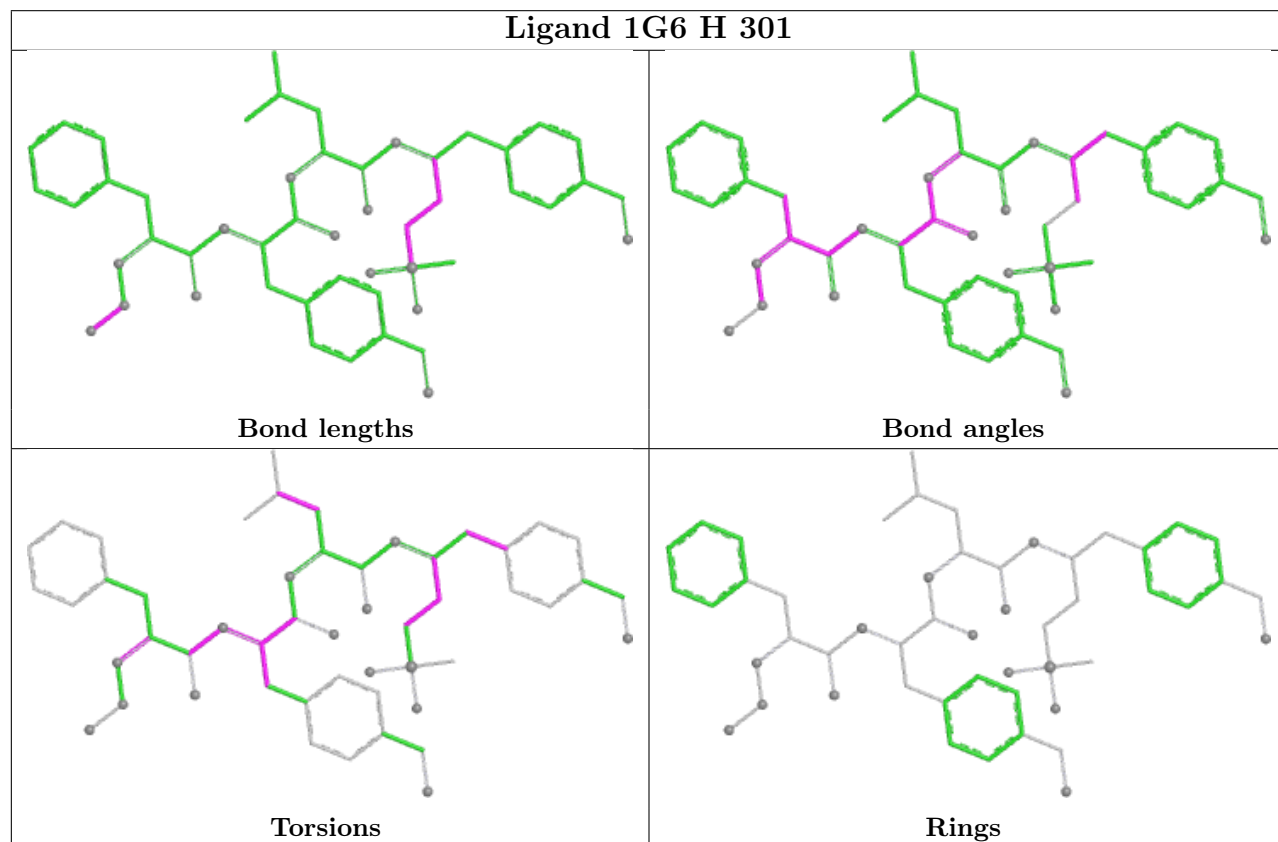
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

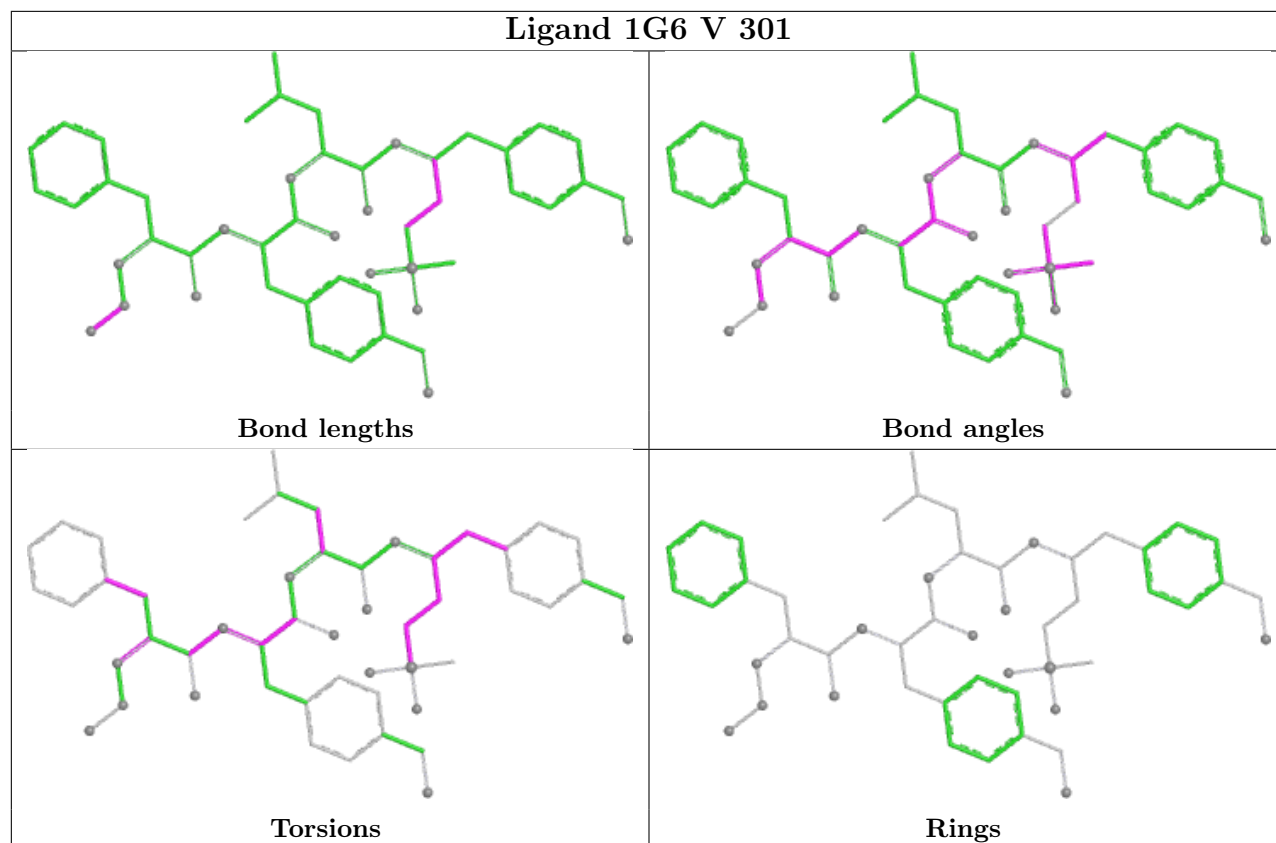


Ligand 1G6 K 301



Ligand 1G6 H 301





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	250/250 (100%)	-0.26	1 (0%) 88 76	72, 91, 117, 134	0
1	O	250/250 (100%)	-0.31	0 100 100	74, 92, 120, 135	0
2	B	244/258 (94%)	-0.16	2 (0%) 82 65	69, 93, 134, 152	0
2	P	244/258 (94%)	-0.09	4 (1%) 70 49	76, 96, 127, 145	0
3	C	241/254 (94%)	-0.19	1 (0%) 88 76	70, 94, 134, 170	0
3	Q	241/254 (94%)	0.03	6 (2%) 58 37	83, 114, 158, 196	0
4	D	242/260 (93%)	-0.18	3 (1%) 76 58	73, 95, 125, 142	0
4	R	242/260 (93%)	-0.09	2 (0%) 82 65	81, 105, 137, 169	0
5	E	233/234 (99%)	-0.13	0 100 100	77, 106, 138, 149	0
5	S	233/234 (99%)	0.00	7 (3%) 52 31	76, 106, 146, 160	0
6	F	244/288 (84%)	-0.22	3 (1%) 76 58	77, 100, 132, 154	0
6	T	244/288 (84%)	-0.30	2 (0%) 82 65	69, 90, 130, 158	0
7	G	243/252 (96%)	-0.18	0 100 100	72, 95, 119, 149	0
7	U	243/252 (96%)	-0.25	1 (0%) 88 76	67, 84, 107, 138	0
8	H	222/232 (95%)	-0.35	0 100 100	66, 79, 102, 120	0
8	V	222/232 (95%)	-0.35	1 (0%) 87 73	61, 76, 95, 126	0
9	I	204/205 (99%)	-0.43	0 100 100	60, 76, 96, 104	0
9	W	204/205 (99%)	-0.39	1 (0%) 87 73	66, 76, 97, 110	0
10	J	198/198 (100%)	-0.28	4 (2%) 65 44	61, 77, 99, 138	0
10	X	198/198 (100%)	-0.24	4 (2%) 65 44	67, 82, 99, 144	0
11	K	212/212 (100%)	-0.06	6 (2%) 55 34	65, 80, 100, 107	0
11	Y	212/212 (100%)	-0.11	4 (1%) 66 45	67, 81, 108, 114	0
12	L	222/222 (100%)	-0.29	1 (0%) 87 73	60, 80, 111, 120	0
12	Z	222/222 (100%)	-0.26	1 (0%) 87 73	64, 83, 111, 121	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	233/233 (100%)	-0.40	1 (0%) 88 76	62, 81, 96, 102	0
13	a	233/233 (100%)	-0.41	2 (0%) 81 63	64, 78, 95, 98	0
14	N	196/196 (100%)	-0.37	0 100 100	65, 74, 95, 103	0
14	b	196/196 (100%)	-0.43	0 100 100	63, 72, 91, 101	0
All	All	6368/6588 (96%)	-0.24	57 (0%) 81 63	60, 87, 130, 196	0

The worst 5 of 57 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
11	Y	209	ASN	5.6
11	Y	104	TYR	5.3
11	K	182	GLU	4.9
11	K	104	TYR	4.4
3	Q	50	LEU	4.3

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.

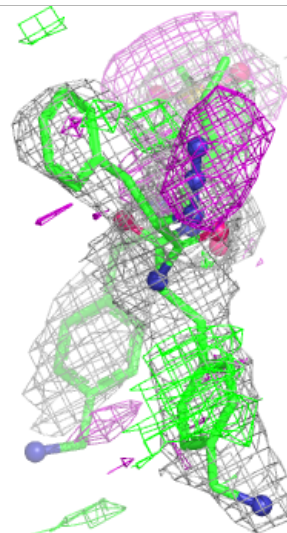
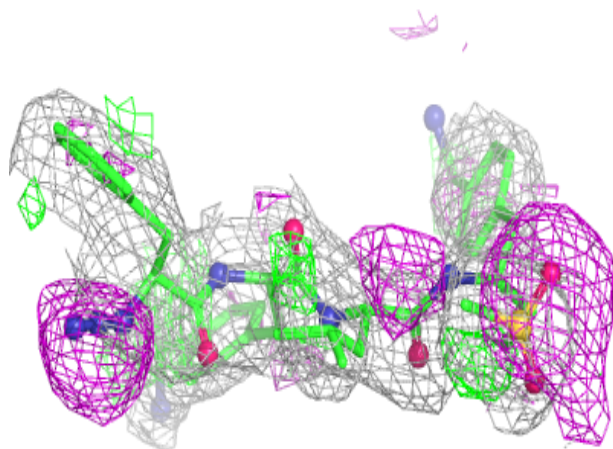
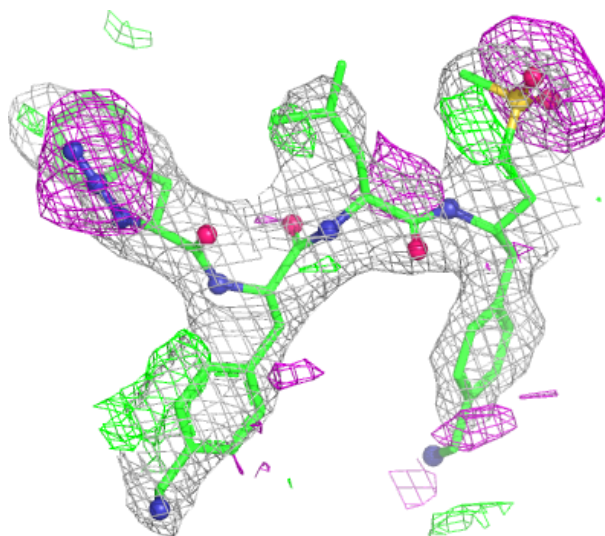
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
15	1G6	V	301	51/51	0.84	0.15	62,68,75,78	0
15	1G6	K	301	51/51	0.87	0.13	58,65,76,79	0
15	1G6	Y	301	51/51	0.89	0.12	57,65,70,71	0
15	1G6	H	301	51/51	0.91	0.11	59,62,67,69	0

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.

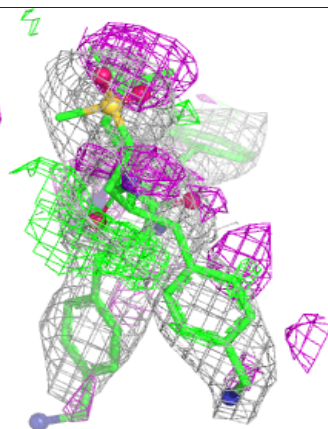
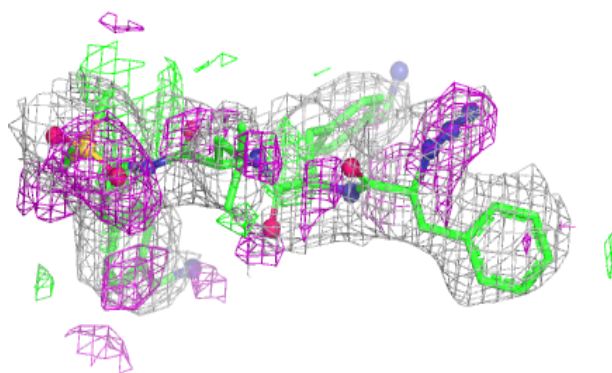
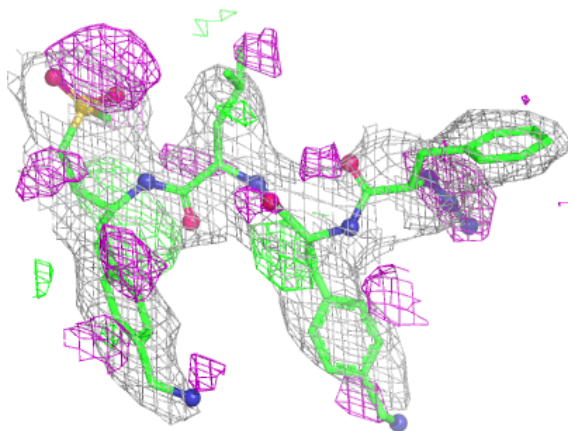
Electron density around 1G6 V 301:

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



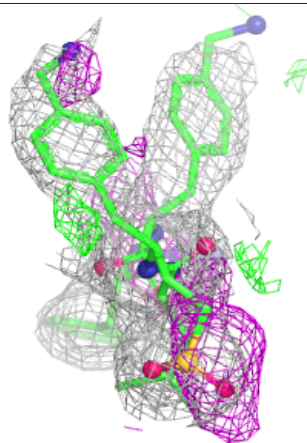
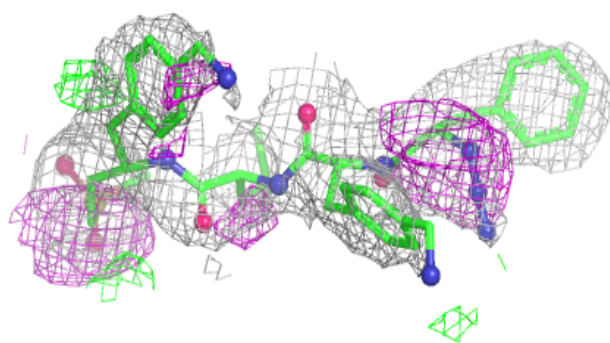
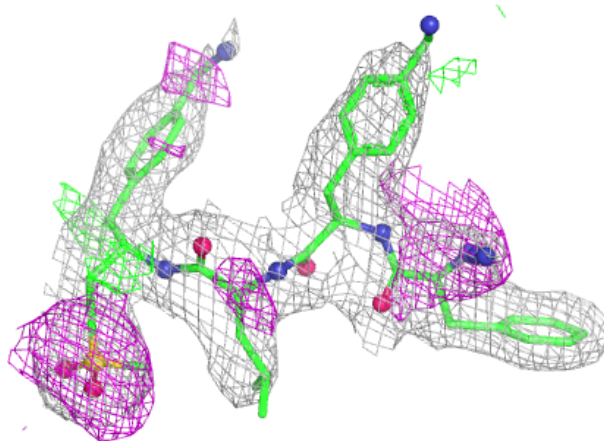
Electron density around 1G6 K 301:

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



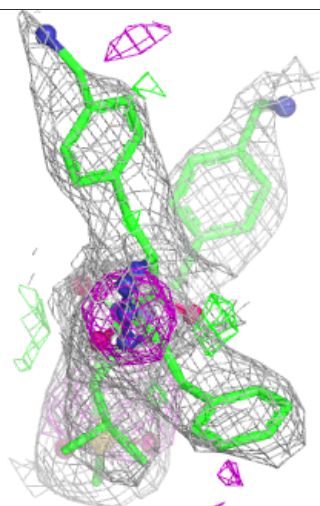
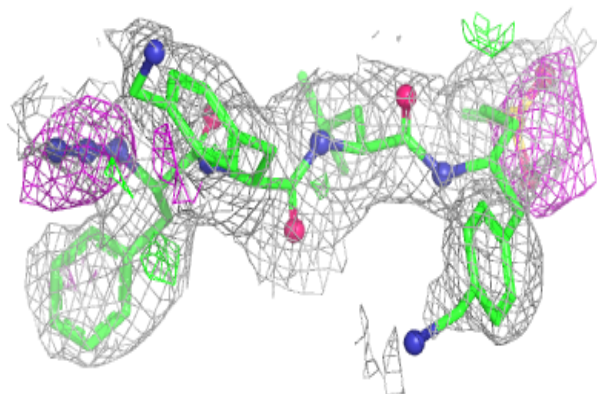
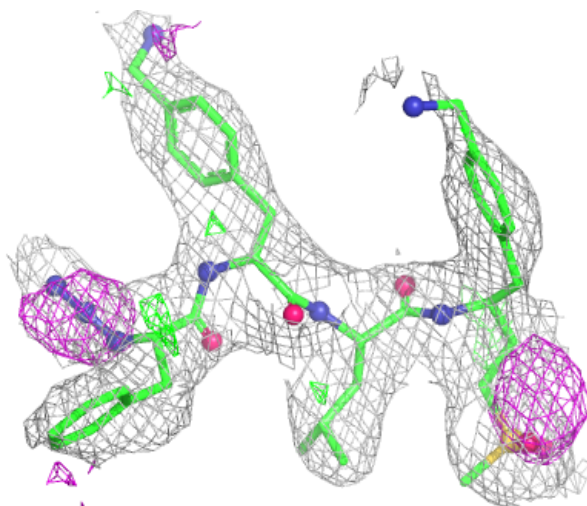
Electron density around 1G6 Y 301:

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



Electron density around 1G6 H 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.