



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 07:44 AM UTC

PDB ID : 8BDJ / pdb_00008bdj
Title : VCB in complex with compound 30
Authors : Sorrell, F.J.; Mueller, J.E.; Lehmann, M.; Wegener, A.
Deposited on : 2022-10-19
Resolution : 2.02 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

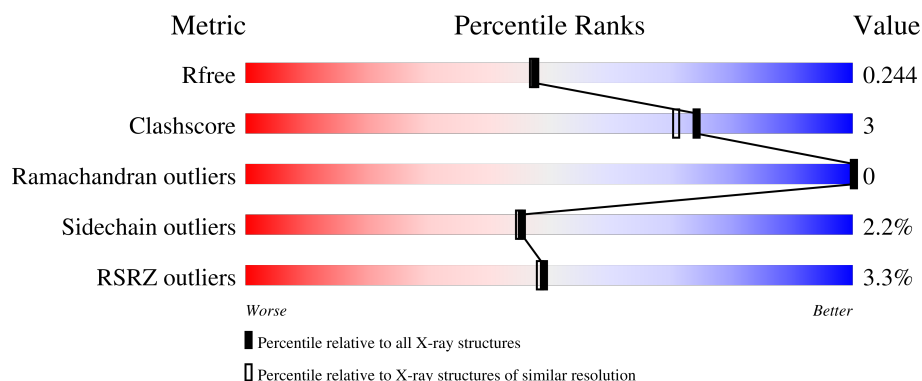
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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	104	0% 94% 5%
1	D	104	5% 93% 5%
1	G	104	2% 95% 5%
2	B	97	8% 80% 11% 7%
2	E	97	2% 85% 5% 7%

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Mol	Chain	Length	Quality of chain
2	H	97	4% 85% 5% 9%
2	K	97	4% 77% 12% 9%
3	C	162	4% 81% 8% 10%
3	F	162	2% 80% 7% 12%
3	I	162	2% 81% 9% 9%
3	L	162	3% 81% 8% 10%
4	J	104	90% 10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11366 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 Elongin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	104	Total	C	N	O	S	0	0	0
			816	516	137	158	5			
1	D	102	Total	C	N	O	S	0	0	0
			802	509	135	153	5			
1	G	104	Total	C	N	O	S	0	0	0
			809	512	135	157	5			

- Molecule 2 is a protein called Elongin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	90	Total	C	N	O	S	0	0	0
			710	457	114	133	6			
2	E	90	Total	C	N	O	S	0	0	0
			715	459	114	135	7			
2	H	88	Total	C	N	O	S	0	0	0
			695	449	111	128	7			
2	K	88	Total	C	N	O	S	0	0	0
			701	452	112	130	7			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	16	MET	-	initiating methionine	UNP Q15369
E	16	MET	-	initiating methionine	UNP Q15369
H	16	MET	-	initiating methionine	UNP Q15369
K	16	MET	-	initiating methionine	UNP Q15369

- Molecule 3 is a protein called von Hippel-Lindau disease tumor suppressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
3	C	145	Total	As	C	N	O	S	0	1	0
			1178	1	750	214	211	2			

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	F	143	Total	As	C	N	O	S	0	0	0
			1160	1	738	207	212	2			
3	I	147	Total	As	C	N	O	S	0	0	0
			1195	1	760	216	216	2			
3	L	146	Total	As	C	N	O	S	0	0	0
			1194	1	758	218	215	2			

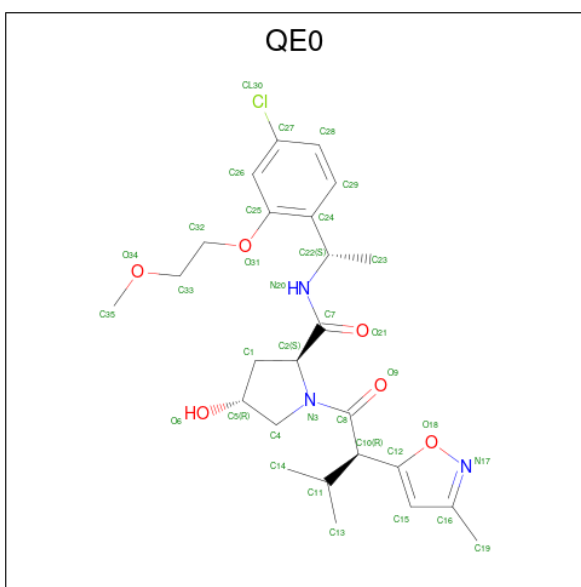
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	52	GLY	-	expression tag	UNP P40337
C	53	SER	-	expression tag	UNP P40337
F	52	GLY	-	expression tag	UNP P40337
F	53	SER	-	expression tag	UNP P40337
I	52	GLY	-	expression tag	UNP P40337
I	53	SER	-	expression tag	UNP P40337
L	52	GLY	-	expression tag	UNP P40337
L	53	SER	-	expression tag	UNP P40337

- Molecule 4 is a protein called Elongin-B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	J	104	Total	As	C	N	O	S	0	0	0
			826	1	522	138	160	5			

- Molecule 5 is (2 {S},4 {R})- {N}-[(1 {S})-1-[4-chloranyl-2-(2-methoxyethoxy)phenyl]ethyl]-1-[(2 {R})-3-methyl-2-(3-methyl-1,2-oxazol-5-yl)butanoyl]-4-oxidanyl-pyrrolidine-2-carboxamide (CCD ID: QE0) (formula: C₂₅H₃₄ClN₃O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total 35	C 25	Cl 1	N 3	O 6	0	0
5	F	1	Total 35	C 25	Cl 1	N 3	O 6	0	0
5	I	1	Total 35	C 25	Cl 1	N 3	O 6	0	0
5	L	1	Total 35	C 25	Cl 1	N 3	O 6	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	49	Total O 49 49	0	0
6	B	31	Total O 31 31	0	0
6	C	50	Total O 50 50	0	0
6	D	20	Total O 20 20	0	0
6	E	17	Total O 17 17	0	0
6	F	43	Total O 43 43	0	0
6	G	26	Total O 26 26	0	0
6	H	22	Total O 22 22	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	I	34	Total 34	O 34	0	0
6	J	47	Total 47	O 47	0	0
6	K	28	Total 28	O 28	0	0
6	L	58	Total 58	O 58	0	0

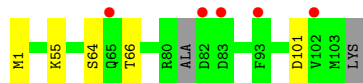
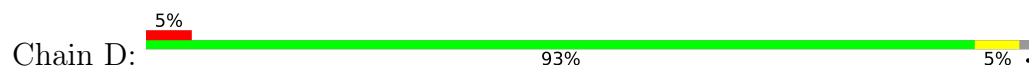
3 Residue-property plots [i](#)

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.

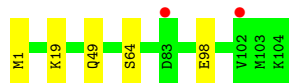
- Molecule 1: Elongin-B



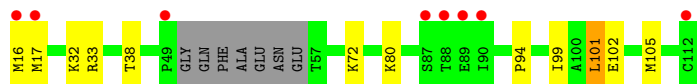
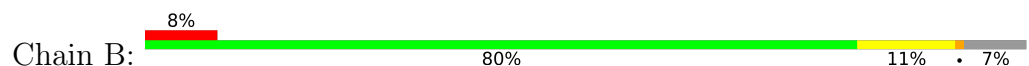
- Molecule 1: Elongin-B



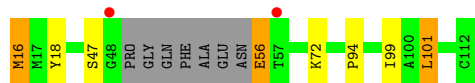
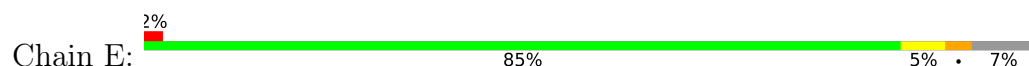
- Molecule 1: Elongin-B



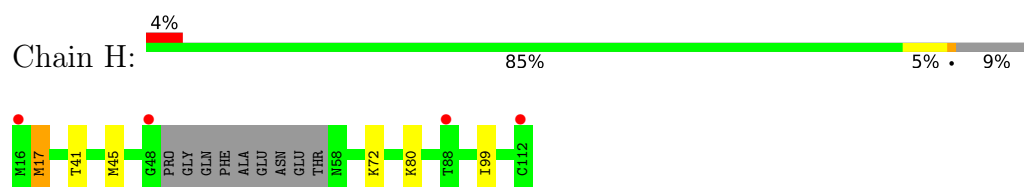
- Molecule 2: Elongin-C



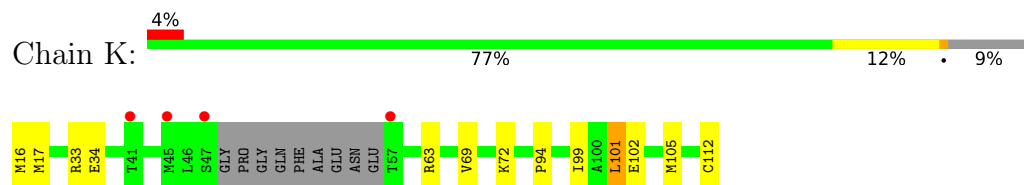
- Molecule 2: Elongin-C



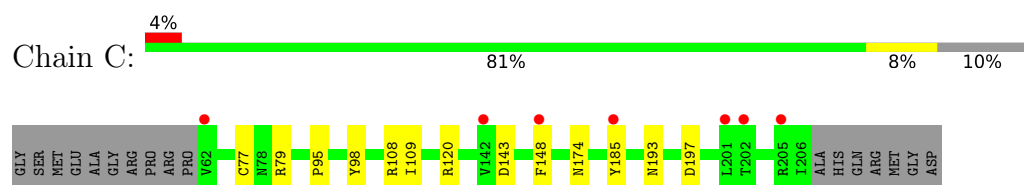
- Molecule 2: Elongin-C



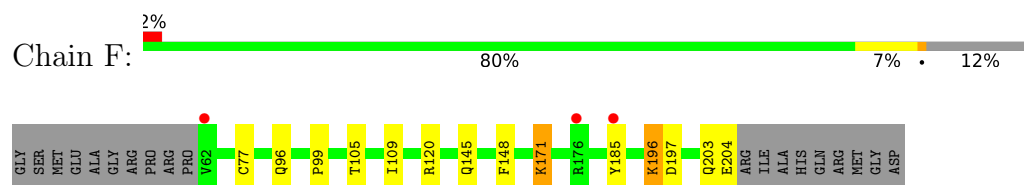
- Molecule 2: Elongin-C



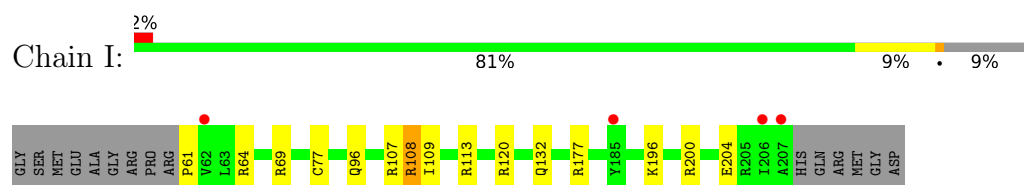
- Molecule 3: von Hippel-Lindau disease tumor suppressor



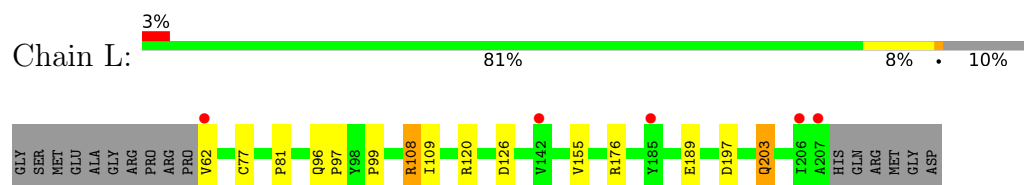
- Molecule 3: von Hippel-Lindau disease tumor suppressor



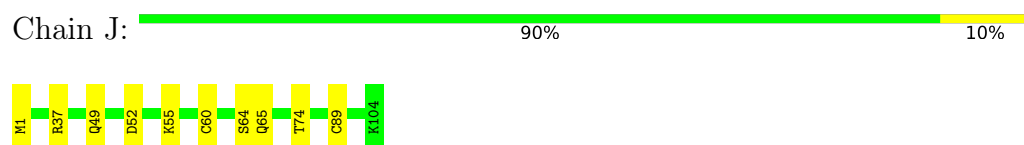
- Molecule 3: von Hippel-Lindau disease tumor suppressor



- Molecule 3: von Hippel-Lindau disease tumor suppressor



- Molecule 4: Elongin-B



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	93.33Å 93.33Å 365.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.33 – 2.02 91.33 – 2.02	Depositor EDS
% Data completeness (in resolution range)	82.2 (91.33-2.02) 82.6 (91.33-2.02)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.02Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.194 , 0.239 0.203 , 0.244	Depositor DCC
R_{free} test set	4449 reflections (4.13%)	wwPDB-VP
Wilson B-factor (Å ²)	44.1	Xtrriage
Anisotropy	0.009	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11366	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.44 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1994e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CAS, QE0

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	1.12	1/832 (0.1%)	1.36	1/1124 (0.1%)
1	D	1.05	0/817	1.31	0/1103
1	G	1.07	0/825	1.30	0/1116
2	B	1.16	0/725	1.30	0/979
2	E	1.12	0/729	1.34	1/982 (0.1%)
2	H	1.13	0/709	1.29	0/955
2	K	1.16	1/715 (0.1%)	1.29	0/964
3	C	1.16	1/1199 (0.1%)	1.34	4/1638 (0.2%)
3	F	1.13	1/1180 (0.1%)	1.37	1/1611 (0.1%)
3	I	1.11	0/1216	1.38	3/1660 (0.2%)
3	L	1.23	4/1214 (0.3%)	1.38	3/1656 (0.2%)
4	J	1.13	1/832 (0.1%)	1.32	0/1121
All	All	1.13	9/10993 (0.1%)	1.34	13/14909 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	95	PRO	C-O	-6.92	1.16	1.23
1	A	34	ILE	C-O	-6.45	1.16	1.23
4	J	74	THR	C-O	6.21	1.31	1.24
3	F	105	THR	C-O	5.92	1.31	1.23
3	L	81	PRO	C-O	-5.72	1.17	1.23
3	L	155	VAL	C-O	-5.63	1.17	1.24
2	K	112	CYS	C-O	5.59	1.34	1.23
3	L	97	PRO	C-O	-5.51	1.17	1.23
3	L	99	PRO	C-O	-5.34	1.17	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	126	ASP	CA-C-N	6.95	127.40	120.31
3	L	126	ASP	C-N-CA	6.95	127.40	120.31
3	I	69	ARG	CG-CD-NE	-5.83	99.19	112.00
3	C	185	TYR	CB-CA-C	-5.79	101.17	110.79
2	E	56	GLU	CB-CA-C	5.78	121.07	110.10
3	F	185	TYR	CB-CA-C	-5.66	101.40	110.79
3	C	79	ARG	CG-CD-NE	-5.53	99.84	112.00
3	C	193	ASN	CB-CA-C	5.35	118.93	109.89
3	L	108	ARG	CG-CD-NE	-5.17	100.64	112.00
3	I	96	GLN	CB-CG-CD	-5.08	103.97	112.60
3	I	108	ARG	CG-CD-NE	-5.06	100.87	112.00
1	A	57	LEU	N-CA-C	-5.06	105.77	111.28
3	C	98	TYR	CB-CA-C	5.04	116.78	109.08

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	816	0	811	4	0
1	D	802	0	801	8	0
1	G	809	0	796	4	0
2	B	710	0	707	8	0
2	E	715	0	713	6	0
2	H	695	0	693	4	0
2	K	701	0	701	8	0
3	C	1178	0	1148	6	0
3	F	1160	0	1137	11	0
3	I	1195	0	1181	6	0
3	L	1194	0	1184	5	0
4	J	826	0	822	5	0
5	C	35	0	0	0	0
5	F	35	0	0	1	0
5	I	35	0	0	0	0
5	L	35	0	0	0	0
6	A	49	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	31	0	0	1	0
6	C	50	0	0	1	0
6	D	20	0	0	0	0
6	E	17	0	0	0	0
6	F	43	0	0	3	0
6	G	26	0	0	2	0
6	H	22	0	0	0	0
6	I	34	0	0	1	0
6	J	47	0	0	0	0
6	K	28	0	0	3	0
6	L	58	0	0	2	0
All	All	11366	0	10694	68	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.

All (68) 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:G:49:GLN:HG3	6:G:205:HOH:O	1.85	0.77
3:C:143:ASP:CB	3:I:177:ARG:NH2	2.49	0.76
3:F:96:GLN:OE1	6:F:401:HOH:O	2.04	0.75
2:B:17:MET:HE3	6:B:219:HOH:O	1.86	0.74
2:K:105:MET:HG2	6:L:457:HOH:O	1.93	0.68
2:K:16:MET:HE2	6:K:221:HOH:O	1.95	0.67
1:A:82:ASP:OD1	1:A:83:ASP:N	2.28	0.66
2:E:56:GLU:O	2:E:56:GLU:HG3	1.99	0.61
3:I:120:ARG:NH2	6:I:401:HOH:O	2.26	0.59
3:F:120:ARG:NH2	6:F:402:HOH:O	2.26	0.59
2:B:16:MET:CB	2:B:32:LYS:HE3	2.32	0.58
3:C:143:ASP:CB	3:I:177:ARG:HH22	2.16	0.58
1:D:55:LYS:HG3	4:J:65:GLN:NE2	2.18	0.57
2:B:33:ARG:NE	1:D:101:ASP:OD2	2.39	0.56
1:G:1:MET:HE2	1:G:64:SER:CB	2.35	0.56
4:J:1:MET:HE2	4:J:64:SER:HB3	1.87	0.55
2:K:33:ARG:NE	6:K:201:HOH:O	2.26	0.54
1:G:1:MET:HE2	1:G:64:SER:HB3	1.90	0.53
1:D:66:THR:HA	4:J:49:GLN:OE1	2.08	0.52
2:H:41:THR:HG22	2:H:45:MET:HE3	1.92	0.51
3:C:120:ARG:NH2	6:C:401:HOH:O	2.27	0.50
3:F:196:LYS:CD	3:F:196:LYS:C	2.85	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:203:GLN:HE21	3:L:203:GLN:HA	1.77	0.49
1:D:1:MET:HE2	1:D:64:SER:CB	2.42	0.48
3:F:99:PRO:HD2	5:F:301:QE0:CL30	2.51	0.48
2:E:72:LYS:NZ	2:E:94:PRO:O	2.49	0.46
3:C:120:ARG:NH1	3:C:197:ASP:OD2	2.43	0.46
1:G:98:GLU:HG2	6:G:223:HOH:O	2.14	0.46
3:L:109:ILE:C	3:L:109:ILE:HD12	2.40	0.46
3:F:196:LYS:C	3:F:196:LYS:HD3	2.40	0.46
1:D:1:MET:HA	1:D:1:MET:HE3	1.98	0.46
2:B:72:LYS:NZ	2:B:94:PRO:O	2.50	0.45
3:F:171:LYS:H	3:F:171:LYS:HG2	1.55	0.45
3:F:109:ILE:C	3:F:109:ILE:HD12	2.42	0.45
3:I:61:PRO:O	3:I:64:ARG:HD3	2.17	0.44
3:F:203:GLN:O	3:F:204:GLU:HG3	2.17	0.44
3:I:109:ILE:C	3:I:109:ILE:HD12	2.42	0.44
3:L:120:ARG:NH1	3:L:197:ASP:OD2	2.46	0.44
3:F:120:ARG:NH1	3:F:197:ASP:OD2	2.47	0.44
3:C:109:ILE:HD12	3:C:109:ILE:C	2.43	0.44
2:B:102:GLU:HA	2:B:105:MET:HE3	1.99	0.43
1:D:66:THR:HG21	4:J:60:CYS:O	2.17	0.43
2:E:16:MET:HE3	2:E:18:TYR:OH	2.18	0.43
1:A:1:MET:HE2	1:A:64:SER:CB	2.48	0.43
3:L:176:ARG:HD3	3:L:189:GLU:OE2	2.19	0.43
2:B:101:LEU:HD12	2:B:101:LEU:HA	1.83	0.43
2:K:72:LYS:NZ	2:K:94:PRO:O	2.51	0.43
2:K:101:LEU:HD12	2:K:101:LEU:HA	1.86	0.43
3:L:120:ARG:NH2	6:L:401:HOH:O	2.33	0.43
1:D:1:MET:HE3	1:D:1:MET:CA	2.48	0.43
3:F:120:ARG:NE	6:F:402:HOH:O	2.47	0.43
2:B:72:LYS:HG3	2:B:99:ILE:CD1	2.49	0.43
1:A:1:MET:HE2	1:A:64:SER:HB3	2.02	0.42
2:E:72:LYS:HG3	2:E:99:ILE:CD1	2.50	0.42
2:H:72:LYS:HG3	2:H:99:ILE:CD1	2.49	0.42
3:I:200:ARG:HG2	3:I:204:GLU:OE2	2.19	0.42
2:K:72:LYS:HG3	2:K:99:ILE:CD1	2.49	0.42
2:E:16:MET:HE2	2:E:16:MET:HB2	1.82	0.42
2:B:38:THR:HG23	2:B:80:LYS:HD3	2.02	0.42
3:F:148:PHE:CZ	2:H:45:MET:HG2	2.55	0.41
1:D:1:MET:HE2	1:D:64:SER:HB2	2.02	0.41
4:J:52:ASP:HB2	4:J:55:LYS:HG2	2.03	0.41
2:K:33:ARG:CD	6:K:201:HOH:O	2.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:69:VAL:HG21	2:K:102:GLU:HB3	2.02	0.41
1:A:82:ASP:OD1	1:A:82:ASP:C	2.64	0.40
2:H:17:MET:HE2	2:H:17:MET:HB2	1.97	0.40
2:E:101:LEU:HD12	2:E:101:LEU:HA	1.83	0.40
3:C:174:ASN:HD22	3:C:174:ASN:HA	1.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	102/104 (98%)	97 (95%)	5 (5%)	0	100	100
1	D	98/104 (94%)	95 (97%)	3 (3%)	0	100	100
1	G	102/104 (98%)	96 (94%)	6 (6%)	0	100	100
2	B	86/97 (89%)	86 (100%)	0	0	100	100
2	E	86/97 (89%)	85 (99%)	1 (1%)	0	100	100
2	H	84/97 (87%)	84 (100%)	0	0	100	100
2	K	84/97 (87%)	84 (100%)	0	0	100	100
3	C	143/162 (88%)	140 (98%)	3 (2%)	0	100	100
3	F	140/162 (86%)	136 (97%)	4 (3%)	0	100	100
3	I	144/162 (89%)	140 (97%)	4 (3%)	0	100	100
3	L	143/162 (88%)	140 (98%)	3 (2%)	0	100	100
4	J	101/104 (97%)	98 (97%)	3 (3%)	0	100	100
All	All	1313/1452 (90%)	1281 (98%)	32 (2%)	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	90/92 (98%)	89 (99%)	1 (1%)	65	68
1	D	89/92 (97%)	89 (100%)	0	100	100
1	G	88/92 (96%)	87 (99%)	1 (1%)	65	68
2	B	80/86 (93%)	79 (99%)	1 (1%)	61	63
2	E	81/86 (94%)	78 (96%)	3 (4%)	30	24
2	H	77/86 (90%)	75 (97%)	2 (3%)	40	38
2	K	79/86 (92%)	75 (95%)	4 (5%)	21	14
3	C	129/147 (88%)	127 (98%)	2 (2%)	55	56
3	F	130/147 (88%)	127 (98%)	3 (2%)	44	43
3	I	134/147 (91%)	129 (96%)	5 (4%)	30	24
3	L	134/147 (91%)	130 (97%)	4 (3%)	36	32
4	J	91/91 (100%)	90 (99%)	1 (1%)	65	68
All	All	1202/1299 (92%)	1175 (98%)	27 (2%)	45	45

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
2	B	101	LEU
3	C	108	ARG
3	C	148	PHE
2	E	16	MET
2	E	47	SER
2	E	101	LEU
3	F	145	GLN
3	F	171	LYS
3	F	196	LYS
1	G	19	LYS
2	H	17	MET
2	H	80	LYS
3	I	107	ARG

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Mol	Chain	Res	Type
3	I	108	ARG
3	I	113	ARG
3	I	132	GLN
3	I	196	LYS
4	J	37	ARG
2	K	17	MET
2	K	34	GLU
2	K	63	ARG
2	K	101	LEU
3	L	62	VAL
3	L	96	GLN
3	L	108	ARG
3	L	203	GLN

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

Mol	Chain	Res	Type
2	B	61	ASN
3	C	145	GLN
3	C	174	ASN
1	D	65	GLN
3	F	145	GLN
3	F	174	ASN
3	F	195	GLN
3	F	203	GLN
3	I	174	ASN
3	I	191	HIS
3	L	132	GLN
3	L	145	GLN
3	L	174	ASN
3	L	203	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

5 non-standard protein/DNA/RNA residues 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$
4	CAS	J	89	4	5,8,9	1.52	1 (20%)	1,9,11	1.33	0
3	CAS	I	77	3	5,8,9	1.29	1 (20%)	1,9,11	1.61	0
3	CAS	F	77	3	5,8,9	1.20	1 (20%)	1,9,11	1.15	0
3	CAS	L	77	3	5,8,9	1.61	1 (20%)	1,9,11	1.34	0
3	CAS	C	77	3	5,8,9	1.97	2 (40%)	1,9,11	1.58	0

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
4	CAS	J	89	4	-	0/0/7/9	-
3	CAS	I	77	3	-	0/0/7/9	-
3	CAS	F	77	3	-	0/0/7/9	-
3	CAS	L	77	3	-	0/0/7/9	-
3	CAS	C	77	3	-	0/0/7/9	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	77	CAS	AS-CE2	3.34	2.04	1.96
3	L	77	CAS	AS-CE2	2.95	2.03	1.96
4	J	89	CAS	AS-CE2	2.65	2.03	1.96
3	C	77	CAS	AS-CE1	2.47	2.02	1.96
3	F	77	CAS	AS-CE2	2.18	2.01	1.96
3	I	77	CAS	AS-CE1	2.06	2.01	1.96

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	$\# Z > 2$
5	QE0	L	301	-	36,37,37	0.48	0	48,52,52	0.59	1 (2%)
5	QE0	C	301	-	36,37,37	0.72	1 (2%)	48,52,52	0.90	2 (4%)
5	QE0	I	301	-	36,37,37	0.37	0	48,52,52	0.65	1 (2%)
5	QE0	F	301	-	36,37,37	0.41	0	48,52,52	0.78	2 (4%)

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	QE0	L	301	-	-	6/33/45/45	0/3/3/3
5	QE0	C	301	-	-	11/33/45/45	0/3/3/3
5	QE0	I	301	-	-	7/33/45/45	0/3/3/3
5	QE0	F	301	-	-	9/33/45/45	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	301	QE0	C16-N17	3.27	1.36	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	QE0	C13-C11-C10	4.10	121.38	111.68
5	F	301	QE0	C13-C11-C10	3.80	120.67	111.68
5	C	301	QE0	C14-C11-C10	2.59	117.81	111.68
5	L	301	QE0	C1-C2-C7	-2.36	106.84	111.36
5	F	301	QE0	C14-C11-C10	2.20	116.88	111.68
5	I	301	QE0	C14-C11-C10	-2.00	106.94	111.68

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	301	QE0	C8-C10-C11-C13
5	C	301	QE0	C8-C10-C11-C14
5	C	301	QE0	C12-C10-C11-C13
5	C	301	QE0	C12-C10-C11-C14
5	F	301	QE0	C8-C10-C11-C13
5	F	301	QE0	C8-C10-C11-C14
5	F	301	QE0	C12-C10-C11-C13
5	F	301	QE0	C12-C10-C11-C14
5	F	301	QE0	C8-C10-C12-O18
5	I	301	QE0	C11-C10-C12-O18
5	L	301	QE0	C11-C10-C12-O18
5	F	301	QE0	O31-C32-C33-O34
5	I	301	QE0	O31-C32-C33-O34
5	I	301	QE0	C33-C32-O31-C25
5	L	301	QE0	C26-C25-O31-C32
5	C	301	QE0	C11-C10-C8-O9
5	F	301	QE0	C26-C25-O31-C32
5	C	301	QE0	C11-C10-C8-N3
5	C	301	QE0	C8-C10-C12-C15
5	C	301	QE0	C8-C10-C12-O18
5	C	301	QE0	C11-C10-C12-O18
5	F	301	QE0	C8-C10-C12-C15
5	I	301	QE0	C8-C10-C12-C15
5	I	301	QE0	C8-C10-C12-O18
5	L	301	QE0	C8-C10-C12-C15
5	L	301	QE0	C8-C10-C12-O18
5	L	301	QE0	C24-C25-O31-C32
5	F	301	QE0	C24-C25-O31-C32
5	C	301	QE0	C33-C32-O31-C25
5	L	301	QE0	C33-C32-O31-C25
5	I	301	QE0	C32-C33-O34-C35
5	I	301	QE0	C26-C25-O31-C32

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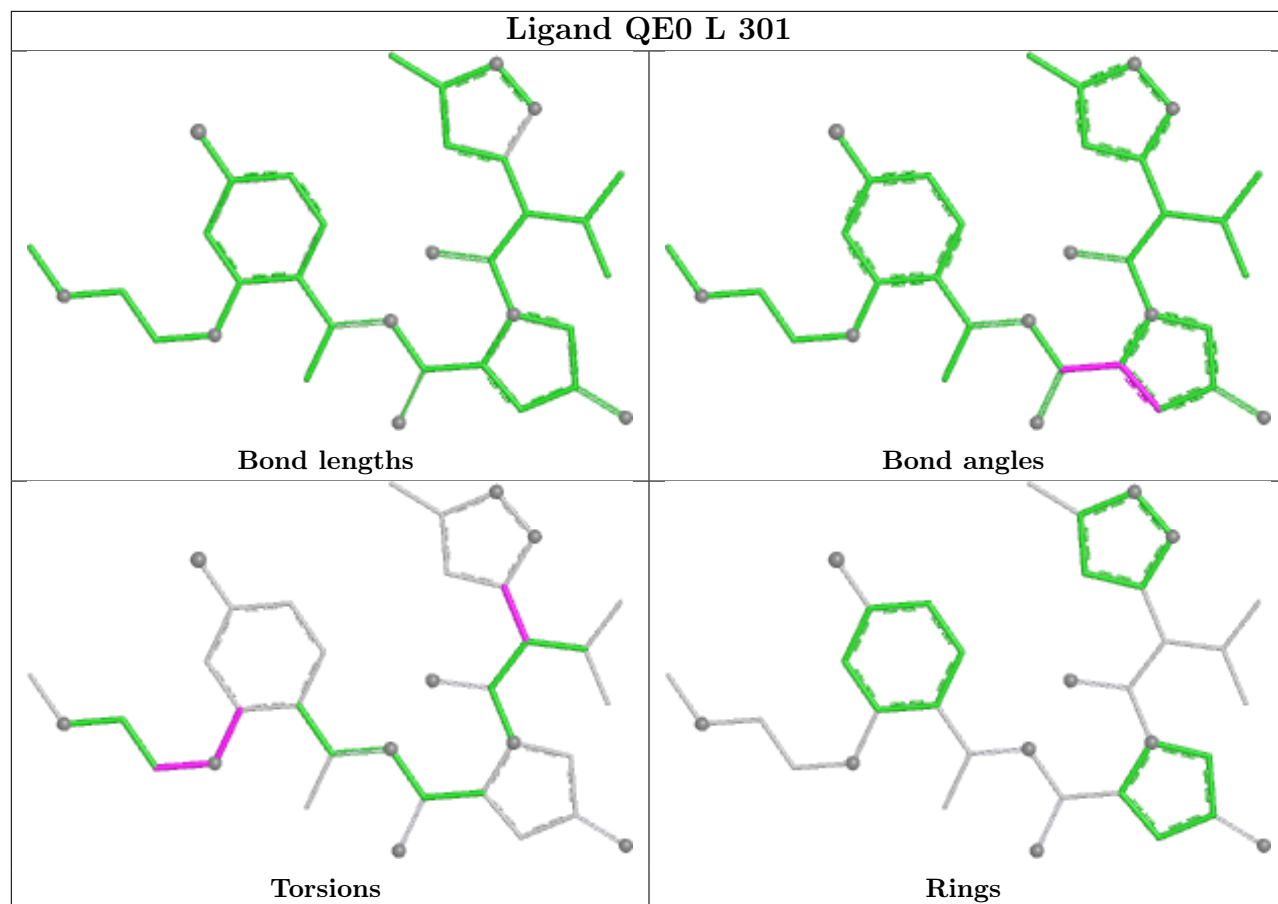
Mol	Chain	Res	Type	Atoms
5	C	301	QE0	C26-C25-O31-C32

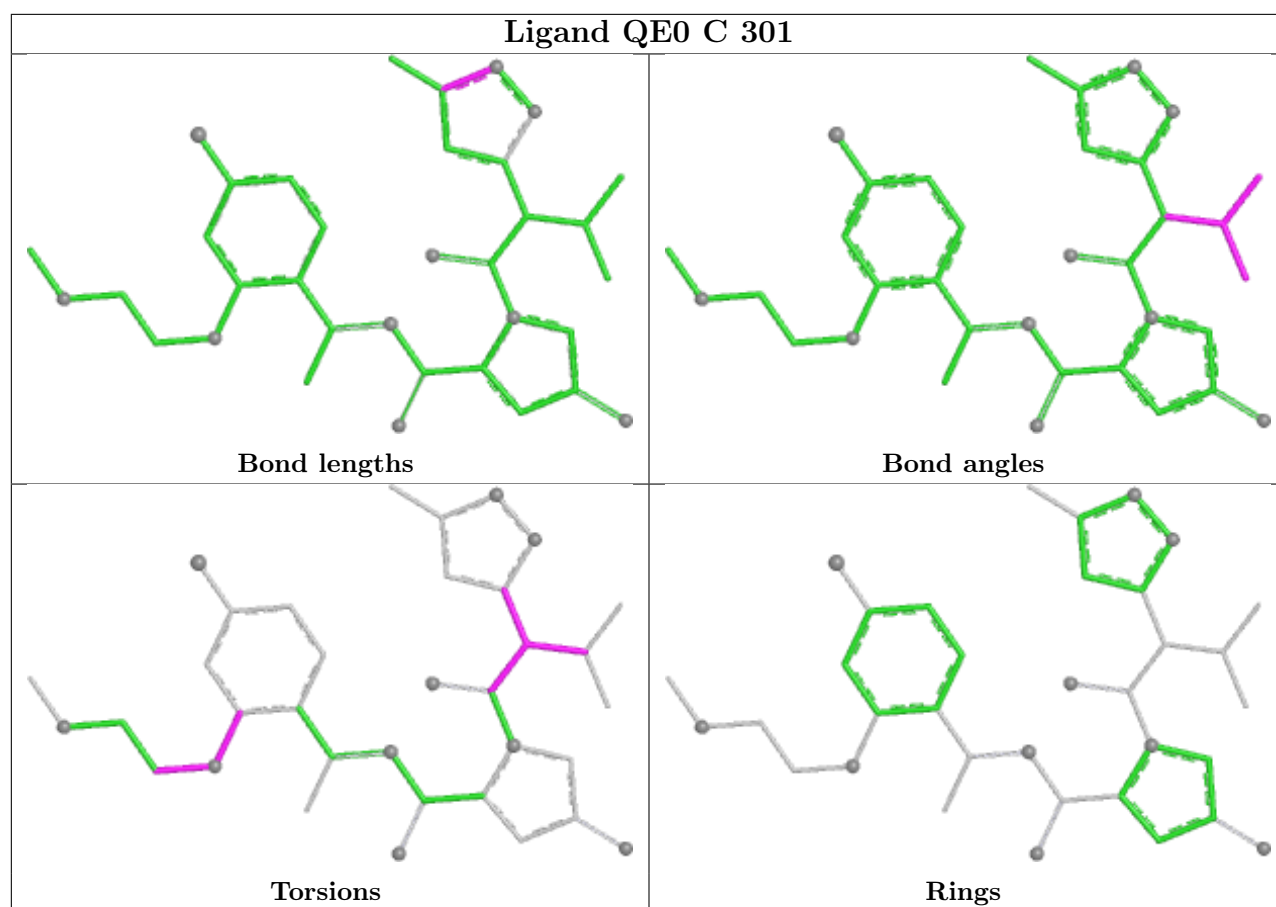
There are no ring outliers.

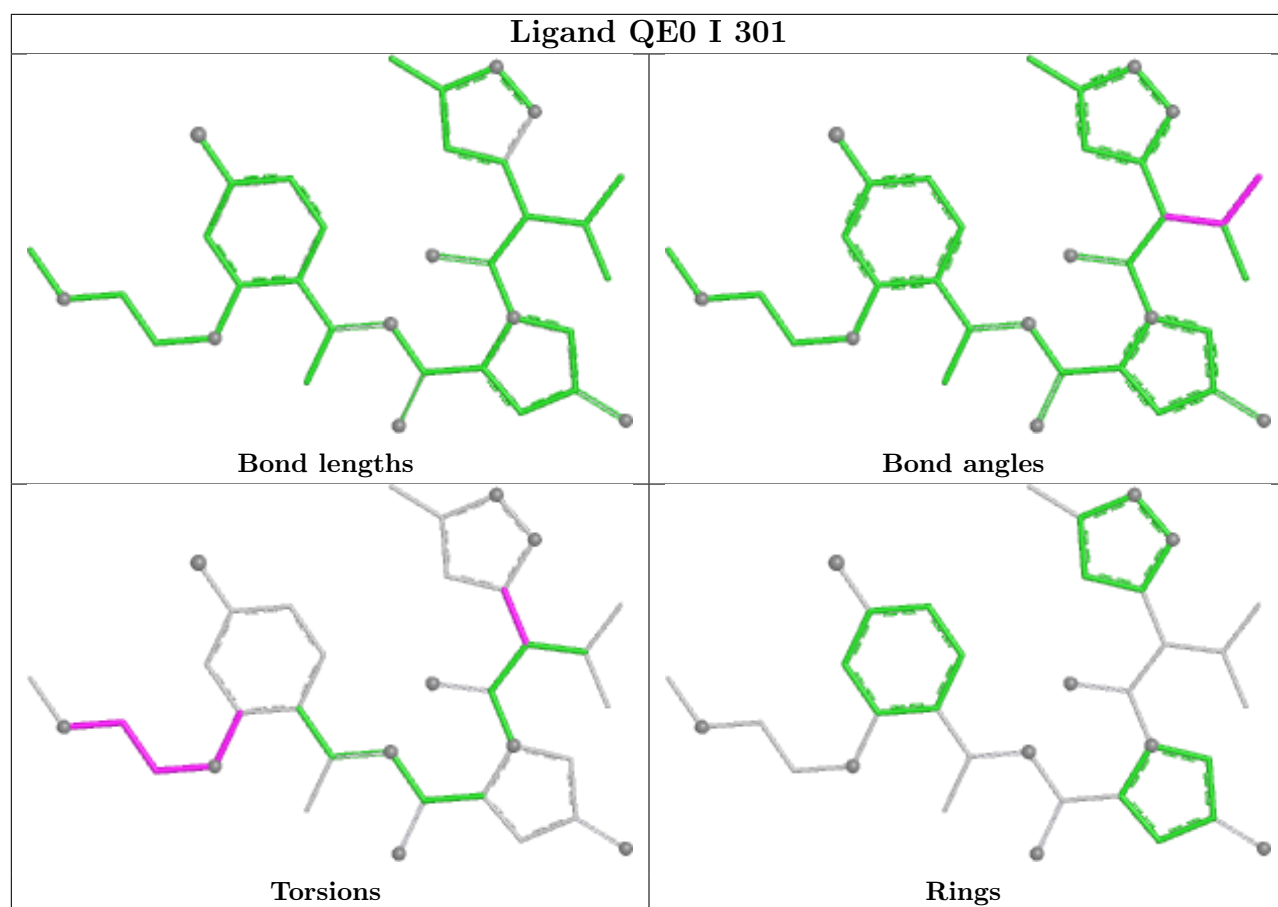
1 monomer is involved in 1 short contact:

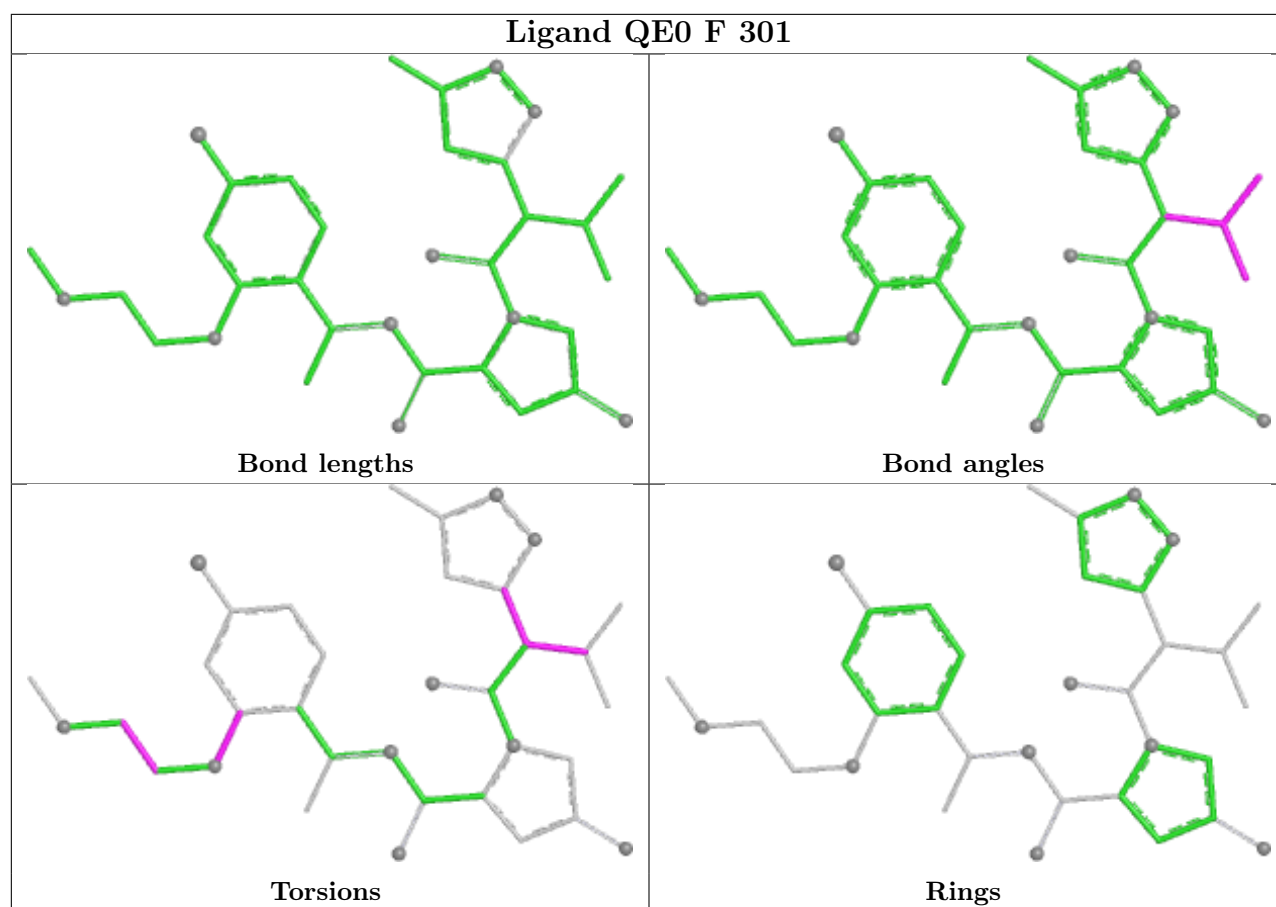
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	301	QE0	1	0

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	104/104 (100%)	-0.10	1 (0%) 79 80	34, 47, 81, 103	0
1	D	102/104 (98%)	0.52	5 (4%) 35 34	44, 69, 102, 106	0
1	G	104/104 (100%)	0.56	2 (1%) 66 66	44, 68, 90, 100	0
2	B	90/97 (92%)	0.17	8 (8%) 15 14	33, 44, 100, 106	0
2	E	90/97 (92%)	0.48	2 (2%) 62 62	41, 60, 95, 106	0
2	H	88/97 (90%)	0.56	4 (4%) 38 37	44, 56, 93, 115	0
2	K	88/97 (90%)	0.27	4 (4%) 38 37	33, 45, 92, 100	0
3	C	144/162 (88%)	0.23	7 (4%) 35 34	29, 49, 96, 137	1 (0%)
3	F	142/162 (87%)	0.19	3 (2%) 63 63	36, 51, 88, 108	0
3	I	146/162 (90%)	0.23	4 (2%) 56 56	38, 52, 91, 135	0
3	L	145/162 (89%)	0.04	5 (3%) 48 47	32, 45, 81, 108	0
4	J	103/104 (99%)	0.00	0 100 100	33, 49, 77, 96	0
All	All	1346/1452 (92%)	0.25	45 (3%) 49 48	29, 52, 94, 137	1 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	90	ILE	5.5
2	H	48	GLY	4.7
3	F	185	TYR	4.4
2	K	47	SER	4.0
3	C	62	VAL	3.7
3	C	185	TYR	3.7
1	D	82	ASP	3.4
1	D	83	ASP	3.3
3	L	62	VAL	3.2
3	L	142	VAL	3.2
3	C	201	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
3	F	62	VAL	2.8
3	C	202	THR	2.7
2	E	48	GLY	2.7
2	B	16	MET	2.7
2	K	45	MET	2.6
2	H	112	CYS	2.6
3	I	185	TYR	2.6
2	B	89	GLU	2.6
2	E	57	THR	2.6
1	D	93	PHE	2.6
3	I	206	ILE	2.6
3	L	185	TYR	2.6
2	B	88	THR	2.4
3	F	176	ARG	2.4
3	I	62	VAL	2.4
2	H	88	THR	2.4
3	L	207	ALA	2.4
1	D	102	VAL	2.4
2	K	57	THR	2.4
3	I	207	ALA	2.3
3	C	148	PHE	2.3
2	B	87	SER	2.3
1	A	83	ASP	2.3
2	B	112	CYS	2.3
3	C	205	ARG	2.3
3	L	206	ILE	2.2
2	K	41	THR	2.2
3	C	142	VAL	2.1
2	H	16	MET	2.1
1	G	83	ASP	2.1
1	D	65	GLN	2.1
1	G	102	VAL	2.1
2	B	17	MET	2.0
2	B	49	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
4	CAS	J	89	9/10	0.83	0.14	52,58,135,144	0
3	CAS	F	77	9/10	0.94	0.12	45,51,84,84	0
3	CAS	C	77	9/10	0.94	0.13	46,51,87,89	0
3	CAS	I	77	9/10	0.95	0.12	41,44,75,76	0
3	CAS	L	77	9/10	0.97	0.10	33,36,64,65	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

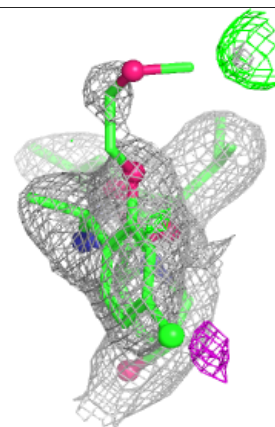
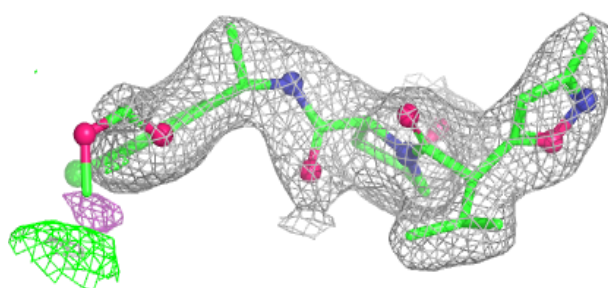
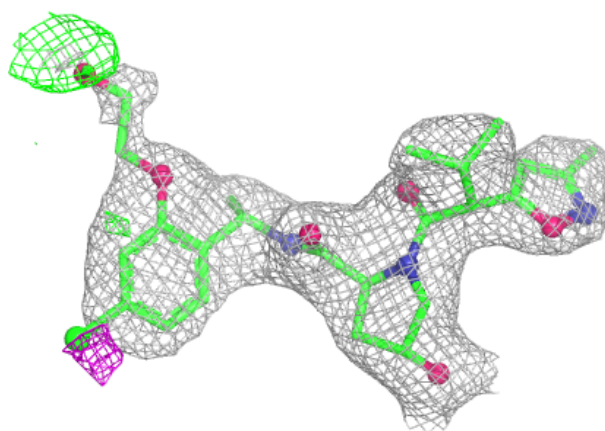
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	QE0	F	301	35/35	0.89	0.13	38,59,104,129	0
5	QE0	L	301	35/35	0.89	0.13	35,50,88,99	0
5	QE0	I	301	35/35	0.90	0.12	37,58,89,93	0
5	QE0	C	301	35/35	0.92	0.12	36,49,84,91	0

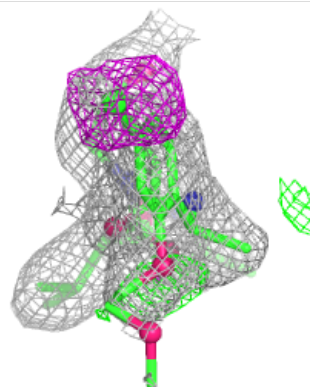
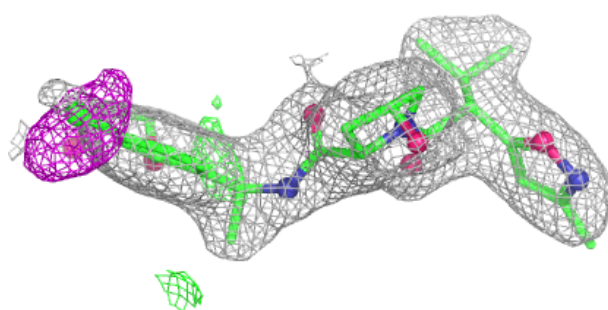
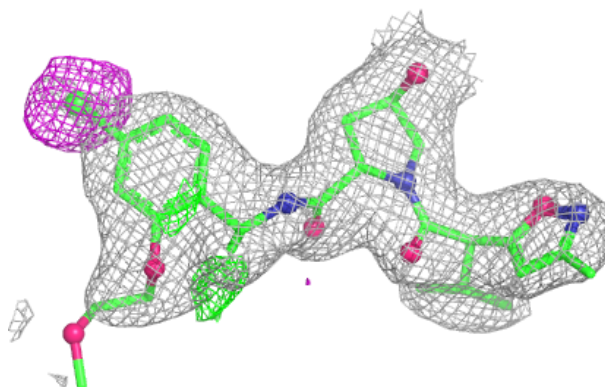
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around QE0 F 301:

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

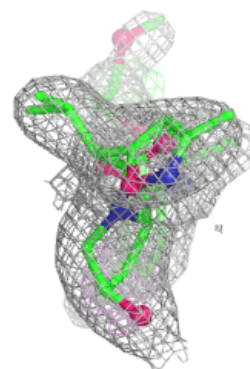
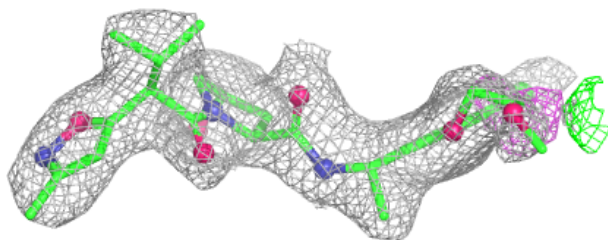
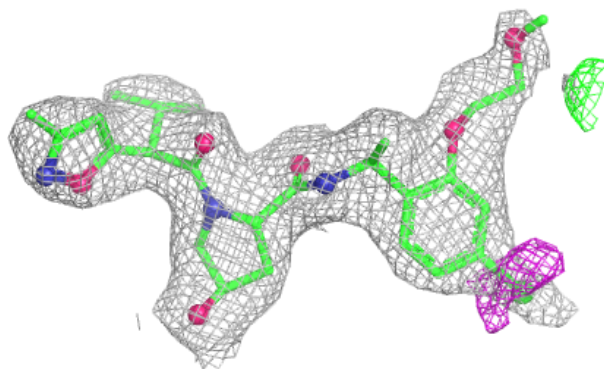
**Electron density around QE0 L 301:**

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



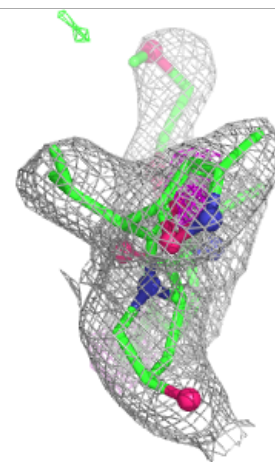
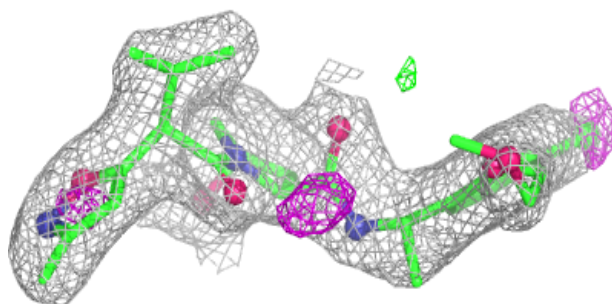
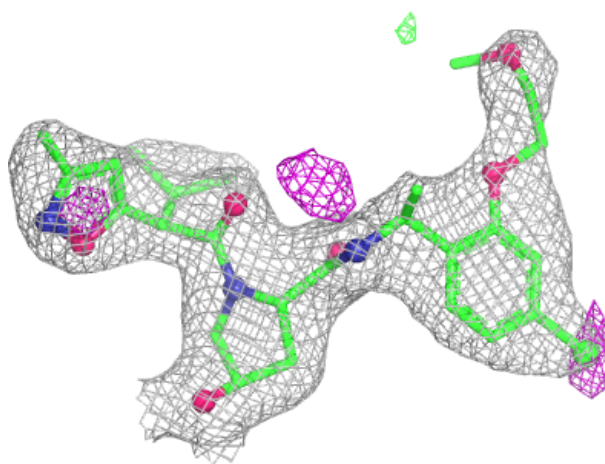
Electron density around QE0 I 301:

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



Electron density around QE0 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.