



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 16, 2020 – 11:45 pm BST

PDB ID : 5TV3  
Title : Crystal structure of the complex of Helicobacter pylori alpha-carbonic anhydrase with (E)-5-(((4-(tert-butyl)phenyl)sulfonyl)imino)-4-methyl-4,5-dihydro-1,3,4-thiadiazole-2-sulfonamide  
Authors : Modak, J.K.; Roujeinikova, A.  
Deposited on : 2016-11-07  
Resolution : 2.90 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

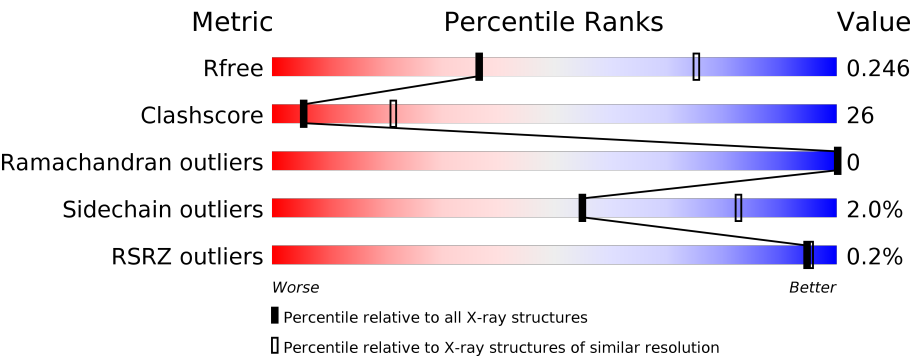
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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 <sub>free</sub>	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	234	<div><div></div><div><div>48%</div><div>47%</div><div>• •</div></div></div>
1	B	234	<div><div></div><div><div>46%</div><div>45%</div><div>• 7%</div></div></div>
1	C	234	<div><div></div><div><div>47%</div><div>49%</div><div>•</div></div></div>
1	D	234	<div><div></div><div><div>39%</div><div>53%</div><div>• 6%</div></div></div>
1	E	234	<div><div></div><div><div>54%</div><div>42%</div><div>•</div></div></div>
1	F	234	<div><div></div><div><div>47%</div><div>47%</div><div>• 5%</div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	234	
1	H	234	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	CL	D	302	-	-	X	-
5	CL	G	302	-	-	X	-

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 14830 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 Alpha-carbonic anhydrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	2	0
			1857	1192	325	336	4			
1	B	218	Total	C	N	O	S	0	1	0
			1787	1147	312	324	4			
1	C	225	Total	C	N	O	S	0	2	0
			1848	1186	323	335	4			
1	D	220	Total	C	N	O	S	0	0	0
			1801	1157	316	324	4			
1	E	226	Total	C	N	O	S	0	1	0
			1853	1188	325	336	4			
1	F	222	Total	C	N	O	S	0	0	0
			1815	1164	318	329	4			
1	G	222	Total	C	N	O	S	0	1	0
			1820	1166	319	331	4			
1	H	225	Total	C	N	O	S	0	0	0
			1841	1180	323	334	4			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

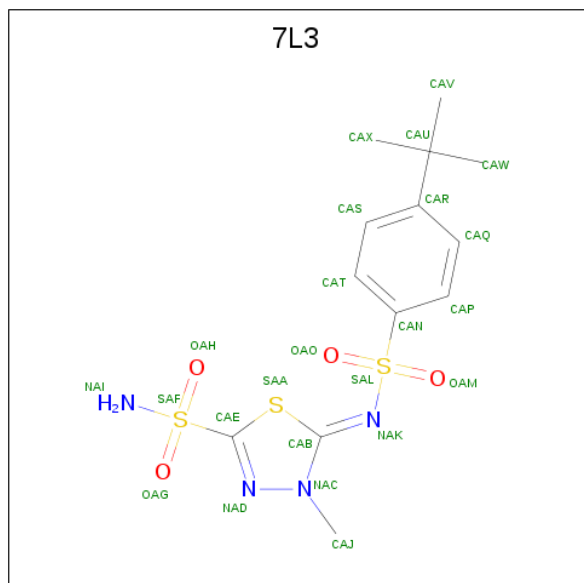
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	H	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

- Molecule 3 is (5Z)-5-[[[(4-tert-butylphenyl)sulfonyl]imino}-4-methyl-4,5-dihydro-1,3,4-thiadiazole-2-sulfonamide (three-letter code: 7L3) (formula: C<sub>13</sub>H<sub>18</sub>N<sub>4</sub>O<sub>4</sub>S<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	B	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	C	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	D	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	E	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	F	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	G	1	Total	C	N	O	S	0	0
			24	13	4	4	3		
3	H	1	Total	C	N	O	S	0	0
			24	13	4	4	3		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

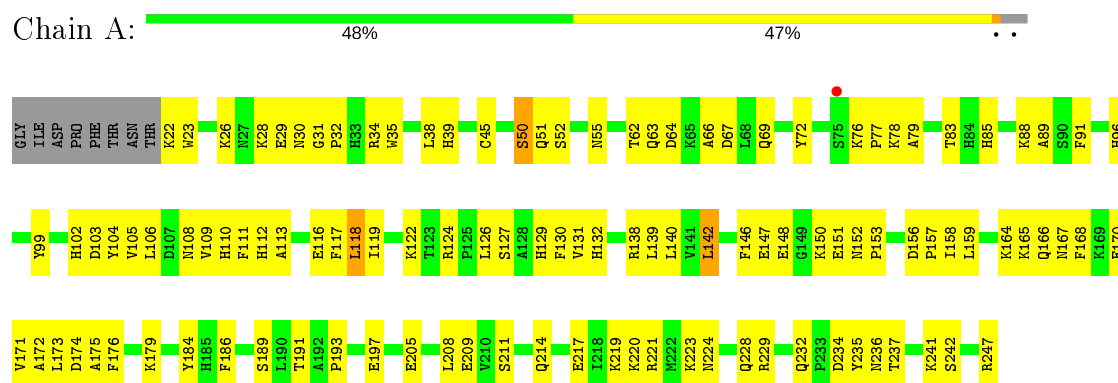
- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		

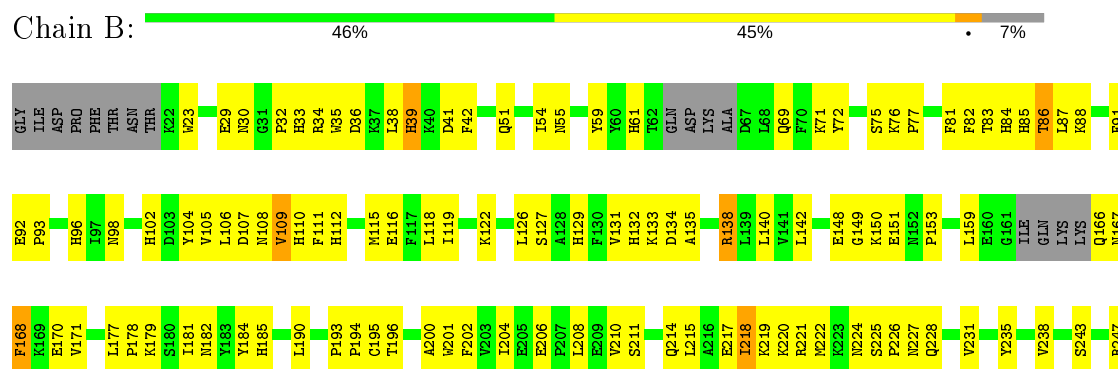
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

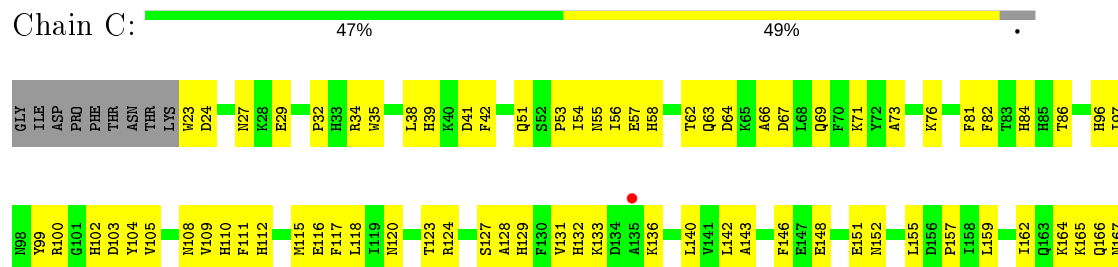
- Molecule 1: Alpha-carbonic anhydrase

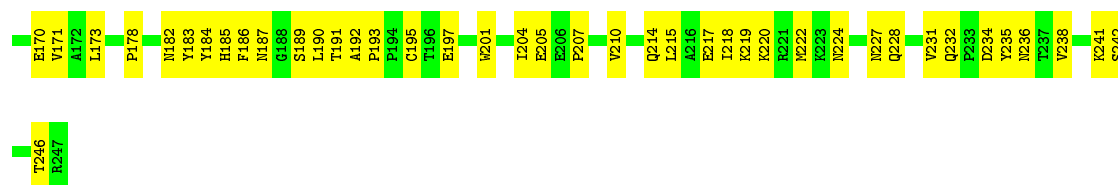


- Molecule 1: Alpha-carbonic anhydrase



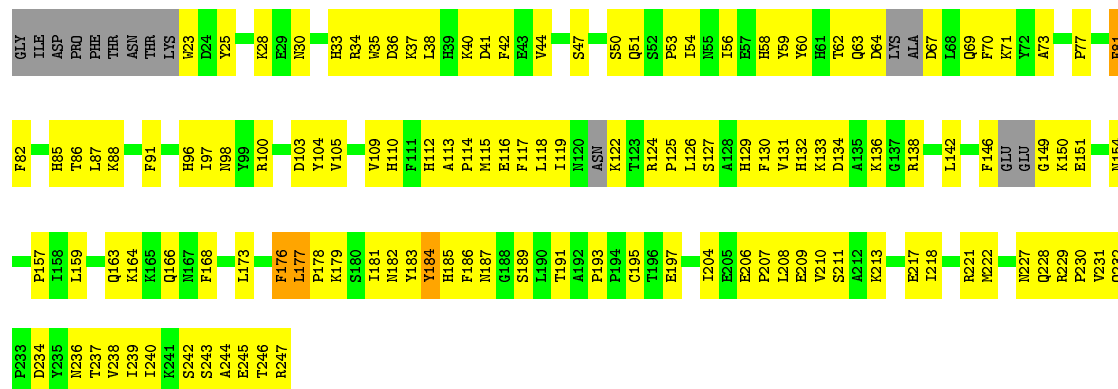
- Molecule 1: Alpha-carbonic anhydrase





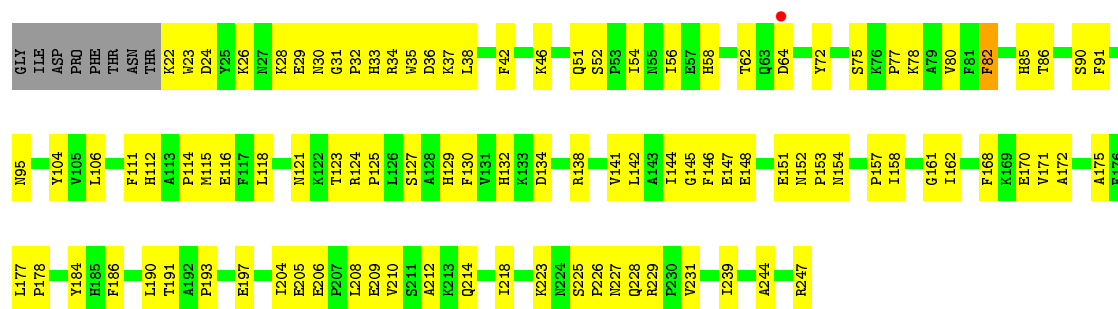
• Molecule 1: Alpha-carbonic anhydrase

Chain D: 39% 53% 6%



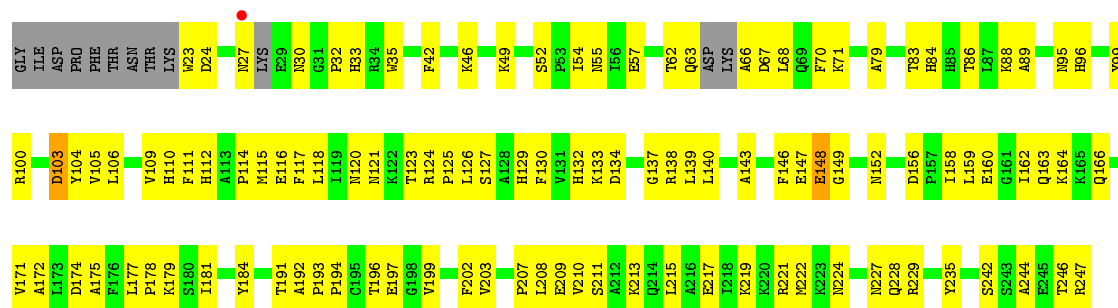
• Molecule 1: Alpha-carbonic anhydrase

Chain E: 54% 42%



• Molecule 1: Alpha-carbonic anhydrase

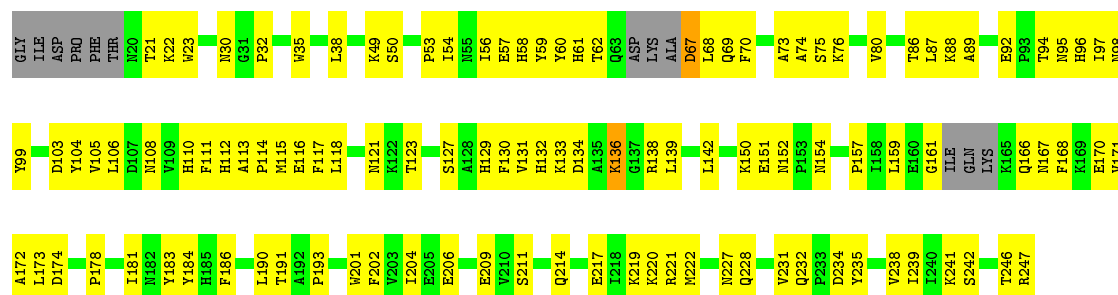
Chain F: 47% 47% 5%





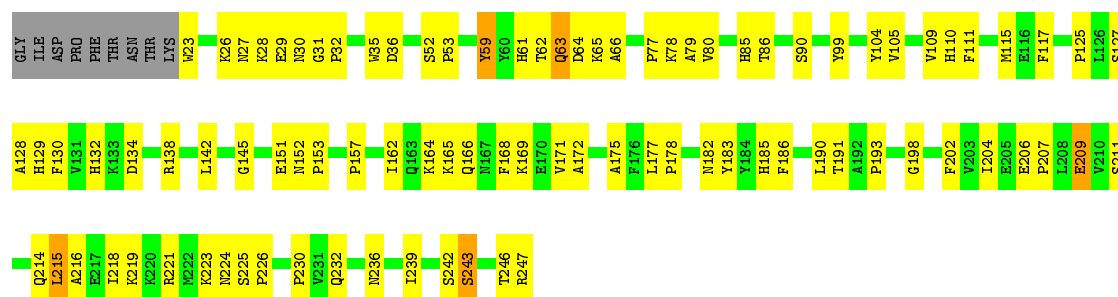
- Molecule 1: Alpha-carbonic anhydrase

Chain G:  47% 47% 5%



- Molecule 1: Alpha-carbonic anhydrase

Chain H:  57% 37%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.56 Å 133.89 Å 166.60 Å 90.00° 90.30° 90.00°	Depositor
Resolution (Å)	29.94 – 2.90 29.94 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.7 (29.94-2.90) 97.2 (29.94-2.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.13 (at 2.90 Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.187 , 0.243 0.183 , 0.246	Depositor DCC
$R_{free}$ test set	2005 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.5	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 24.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.447 for h,-k,-l	Xtriage
Reported twinning fraction	0.500 for h,-k,-l	Depositor
Outliers	1 of 39903 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	14830	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, 7L3, CL

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	0.54	1/1918 (0.1%)	0.80	2/2594 (0.1%)
1	B	0.53	0/1843	0.76	0/2492
1	C	0.56	0/1909	0.79	0/2583
1	D	0.53	0/1853	0.75	1/2503 (0.0%)
1	E	0.58	0/1911	0.80	1/2584 (0.0%)
1	F	0.50	0/1868	0.76	0/2526
1	G	0.52	0/1876	0.76	0/2536
1	H	0.55	0/1896	0.76	1/2565 (0.0%)
All	All	0.54	1/15074 (0.0%)	0.77	5/20383 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	45	CYS	CB-SG	-5.55	1.72	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	LEU	CA-CB-CG	6.92	131.22	115.30
1	E	62	THR	N-CA-C	-6.50	93.45	111.00
1	D	177	LEU	CA-CB-CG	6.03	129.17	115.30
1	H	215	LEU	CA-CB-CG	5.61	128.20	115.30
1	A	118	LEU	CA-CB-CG	-5.52	102.61	115.30

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	1857	0	1825	100	0
1	B	1787	0	1739	91	1
1	C	1848	0	1812	100	0
1	D	1801	0	1759	121	0
1	E	1853	0	1816	83	0
1	F	1815	0	1766	94	0
1	G	1820	0	1773	106	0
1	H	1841	0	1798	76	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
3	A	24	0	0	3	0
3	B	24	0	0	3	0
3	C	24	0	0	3	0
3	D	24	0	0	3	0
3	E	24	0	0	2	0
3	F	24	0	0	4	0
3	G	24	0	0	2	0
3	H	24	0	0	2	0
4	C	6	0	8	1	0
5	D	1	0	0	2	0
5	G	1	0	0	2	0
All	All	14830	0	14296	747	1

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 26.

All (747) 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:D:177:LEU:HD12	1:D:246:THR:CG2	1.68	1.21
1:D:177:LEU:HD12	1:D:246:THR:HG21	1.29	1.11
1:D:177:LEU:HD12	1:D:246:THR:HG23	1.47	0.97
1:A:102:HIS:HB3	1:A:138:ARG:HH22	1.31	0.96
1:B:41:ASP:HA	1:D:40:LYS:HE3	1.50	0.94
1:D:177:LEU:CD1	1:D:246:THR:HG21	2.00	0.89
1:B:72:TYR:OH	1:B:132:HIS:NE2	2.06	0.88
1:D:177:LEU:HB2	1:D:246:THR:HG21	1.54	0.88
1:H:130:PHE:HB2	1:H:142:LEU:HB3	1.57	0.86
1:F:23:TRP:HA	1:F:30:ASN:HB3	1.56	0.86
1:E:115:MET:O	1:E:228:GLN:NE2	2.09	0.84
1:B:87:LEU:HB3	1:B:222:MET:HE2	1.60	0.83
1:C:71:LYS:NZ	1:C:103:ASP:OD2	2.12	0.83
1:C:219:LYS:O	1:C:224:ASN:N	2.10	0.83
1:G:157:PRO:HB2	1:G:171:VAL:HG22	1.61	0.83
1:A:219:LYS:O	1:A:224:ASN:N	2.11	0.82
1:B:177:LEU:O	1:B:247:ARG:NH1	2.12	0.82
1:E:191:THR:OG1	3:E:302:7L3:NAI	2.12	0.81
1:C:159:LEU:HD21	1:C:218:ILE:HG22	1.62	0.81
1:D:54:ILE:HA	1:D:231:VAL:HG13	1.61	0.80
1:G:115:MET:HG3	1:G:127:SER:HB3	1.63	0.80
1:D:149:GLY:N	1:D:210:VAL:O	2.15	0.79
1:B:71:LYS:HB2	1:B:96:HIS:H	1.47	0.78
1:F:104:TYR:HB3	1:F:132:HIS:HB3	1.65	0.78
1:G:136:LYS:HE3	1:G:138:ARG:HG3	1.66	0.77
1:C:32:PRO:HA	1:C:35:TRP:CG	2.19	0.77
1:E:157:PRO:HB2	1:E:171:VAL:HG22	1.66	0.77
1:G:118:LEU:HD23	1:G:121:ASN:HA	1.66	0.77
1:G:76:LYS:HB2	1:G:170:GLU:HG2	1.66	0.77
1:B:33:HIS:CD2	1:B:34:ARG:HG3	2.20	0.77
1:H:151:GLU:HG2	1:H:211:SER:HB3	1.66	0.77
1:D:204:ILE:HG22	1:D:206:GLU:H	1.49	0.76
1:D:23:TRP:HA	1:D:30:ASN:HB3	1.66	0.76
1:C:191:THR:OG1	3:C:303:7L3:NAI	2.18	0.76
1:A:119:ILE:HD13	1:A:205:GLU:HA	1.67	0.75
1:C:71:LYS:HD2	1:C:96:HIS:HB2	1.68	0.75
1:F:219:LYS:O	1:F:224:ASN:N	2.17	0.74
1:D:150:LYS:H	1:D:209:GLU:HG2	1.52	0.74
1:A:152:ASN:ND2	1:A:209:GLU:O	2.21	0.74
1:H:204:ILE:HG22	1:H:206:GLU:H	1.53	0.73
1:H:62:THR:OG1	1:H:63:GLN:N	2.19	0.73
1:D:36:ASP:OD2	1:D:37:LYS:NZ	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:TRP:HA	1:E:30:ASN:HB3	1.70	0.73
1:E:172:ALA:HB1	1:E:175:ALA:HB3	1.69	0.73
1:B:75:SER:O	1:B:171:VAL:N	2.20	0.73
1:C:178:PRO:O	1:C:246:THR:OG1	2.05	0.73
1:E:204:ILE:HG22	1:E:206:GLU:H	1.53	0.73
1:B:102:HIS:HB3	1:B:134:ASP:OD1	1.89	0.72
1:C:24:ASP:O	1:C:35:TRP:NE1	2.19	0.72
1:G:178:PRO:HG2	1:G:206:GLU:HG3	1.71	0.71
1:F:156:ASP:O	1:F:160:GLU:N	2.18	0.71
1:G:54:ILE:HA	1:G:231:VAL:HG13	1.71	0.71
1:D:28:LYS:O	1:D:33:HIS:NE2	2.24	0.71
1:C:115:MET:HG3	1:C:127:SER:HB3	1.70	0.71
1:F:160:GLU:HA	1:F:163:GLN:HB2	1.73	0.71
1:F:181:ILE:HD11	1:F:244:ALA:O	1.91	0.70
1:F:217:GLU:O	1:F:221:ARG:N	2.24	0.70
1:B:82:PHE:HA	1:B:87:LEU:HA	1.72	0.70
1:F:139:LEU:HB2	1:F:199:VAL:HG22	1.74	0.70
1:H:214:GLN:O	1:H:218:ILE:HG13	1.90	0.70
1:A:148:GLU:O	1:F:179:LYS:NZ	2.25	0.70
1:D:130:PHE:HB3	1:D:132:HIS:HE1	1.56	0.69
1:D:230:PRO:O	1:D:232:GLN:NE2	2.24	0.69
1:C:54:ILE:HA	1:C:231:VAL:HG13	1.75	0.69
1:G:234:ASP:O	1:H:138:ARG:NH2	2.25	0.69
1:A:39:HIS:ND1	1:C:41:ASP:OD2	2.23	0.69
1:H:177:LEU:O	1:H:247:ARG:NH1	2.26	0.69
1:C:118:LEU:HD22	1:C:123:THR:HA	1.75	0.68
1:C:104:TYR:HB3	1:C:132:HIS:HB3	1.74	0.68
1:B:134:ASP:CG	1:B:135:ALA:H	1.97	0.68
1:F:184:TYR:N	1:F:203:VAL:O	2.22	0.68
1:B:23:TRP:O	1:B:85:HIS:NE2	2.27	0.67
1:D:71:LYS:HB2	1:D:96:HIS:HB2	1.76	0.67
1:D:62:THR:OG1	1:D:63:GLN:N	2.27	0.67
1:E:82:PHE:CE2	1:E:85:HIS:HA	2.28	0.67
1:G:74:ALA:HA	1:G:172:ALA:HA	1.76	0.67
1:A:72:TYR:OH	1:A:132:HIS:NE2	2.19	0.67
1:H:219:LYS:O	1:H:224:ASN:N	2.17	0.66
1:D:33:HIS:ND1	1:D:34:ARG:HG3	2.10	0.66
1:D:159:LEU:HD21	1:D:218:ILE:HG12	1.78	0.66
1:G:73:ALA:H	1:G:95:ASN:ND2	1.93	0.66
1:B:96:HIS:CD2	1:B:105:VAL:HG22	2.30	0.66
1:A:174:ASP:O	1:A:247:ARG:HD3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:181:ILE:O	1:D:244:ALA:N	2.26	0.66
1:H:29:GLU:HG3	1:H:30:ASN:H	1.60	0.66
1:A:62:THR:OG1	1:A:63:GLN:OE1	2.14	0.65
1:A:116:GLU:HG2	1:A:228:GLN:HG3	1.78	0.65
1:E:54:ILE:HA	1:E:231:VAL:HG13	1.76	0.65
1:E:178:PRO:HD3	1:E:208:LEU:HD21	1.78	0.65
1:D:177:LEU:CB	1:D:246:THR:HG21	2.25	0.65
1:G:73:ALA:H	1:G:95:ASN:HD22	1.44	0.65
1:C:66:ALA:O	1:C:69:GLN:NE2	2.28	0.65
1:E:177:LEU:O	1:E:247:ARG:NH1	2.29	0.65
1:G:59:TYR:HB3	1:G:239:ILE:HB	1.79	0.65
1:C:186:PHE:HA	1:D:238:VAL:HG21	1.78	0.65
1:C:29:GLU:HA	1:C:34:ARG:HE	1.63	0.64
1:F:110:HIS:CE1	1:F:129:HIS:HB2	2.32	0.64
1:F:191:THR:OG1	3:F:302:7L3:NAI	2.31	0.64
1:G:69:GLN:HG3	1:G:70:PHE:N	2.12	0.64
1:A:28:LYS:O	1:A:34:ARG:NH1	2.30	0.64
1:D:177:LEU:CD1	1:D:246:THR:CG2	2.58	0.64
1:G:57:GLU:HG3	1:G:58:HIS:ND1	2.13	0.64
1:A:156:ASP:OD1	1:A:214:GLN:NE2	2.30	0.64
1:A:76:LYS:HG2	1:A:170:GLU:HG2	1.80	0.64
1:C:51:GLN:HB3	1:C:232:GLN:HG3	1.80	0.64
1:A:23:TRP:O	1:A:85:HIS:NE2	2.31	0.64
1:B:185:HIS:HA	1:B:201:TRP:O	1.97	0.64
1:A:96:HIS:ND1	1:A:105:VAL:HG22	2.13	0.63
1:G:150:LYS:HG3	1:G:151:GLU:H	1.63	0.63
1:A:151:GLU:HA	1:A:211:SER:HB3	1.81	0.63
1:C:96:HIS:ND1	1:C:105:VAL:HG22	2.14	0.63
1:F:49:LYS:N	1:F:197:GLU:OE2	2.31	0.63
1:A:130:PHE:HB2	1:A:142:LEU:HB3	1.79	0.63
1:G:32:PRO:HA	1:G:35:TRP:CD2	2.34	0.63
1:C:64:ASP:H	1:H:65:LYS:HB2	1.62	0.63
1:A:119:ILE:HG21	1:A:205:GLU:HG3	1.79	0.63
1:H:157:PRO:HB2	1:H:171:VAL:HG22	1.79	0.63
1:A:124:ARG:NH2	1:A:205:GLU:O	2.32	0.63
1:C:55:ASN:ND2	1:C:57:GLU:OE1	2.28	0.63
1:G:112:HIS:HA	1:G:227:ASN:ND2	2.14	0.63
1:B:30:ASN:HA	1:B:38:LEU:HD11	1.80	0.62
1:E:130:PHE:HB2	1:E:142:LEU:HB3	1.79	0.62
1:F:156:ASP:HA	1:F:159:LEU:HB2	1.79	0.62
1:F:33:HIS:O	1:F:46:LYS:HE2	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:PHE:HB2	1:B:222:MET:HE3	1.80	0.62
1:F:146:PHE:CE1	1:F:208:LEU:HB2	2.34	0.62
1:F:177:LEU:HB2	1:F:247:ARG:HH22	1.64	0.62
1:E:32:PRO:HA	1:E:35:TRP:CG	2.35	0.62
1:G:159:LEU:HD11	1:G:217:GLU:HG2	1.81	0.62
1:E:104:TYR:HB3	1:E:132:HIS:HB3	1.81	0.62
1:G:112:HIS:HA	1:G:227:ASN:HD21	1.64	0.62
1:E:24:ASP:OD1	1:E:30:ASN:HB2	2.00	0.62
1:F:118:LEU:HD23	1:F:121:ASN:HA	1.81	0.62
1:G:53:PRO:HG3	1:G:116:GLU:HB3	1.79	0.62
1:B:194:PRO:HG2	1:B:196:THR:OG1	2.00	0.62
1:E:35:TRP:CH2	1:E:193:PRO:HD3	2.35	0.62
1:B:54:ILE:HA	1:B:231:VAL:HG13	1.81	0.61
1:H:152:ASN:HB2	1:H:209:GLU:O	1.99	0.61
1:D:62:THR:OG1	1:D:63:GLN:OE1	2.19	0.61
1:G:134:ASP:OD1	1:G:138:ARG:N	2.33	0.61
1:E:86:THR:OG1	1:E:111:PHE:O	2.18	0.61
1:C:124:ARG:NH1	1:C:207:PRO:HB3	2.16	0.61
1:C:152:ASN:HB3	1:C:155:LEU:HG	1.83	0.61
1:D:113:ALA:HB2	1:D:126:LEU:HD13	1.83	0.61
1:A:191:THR:OG1	3:A:302:7L3:NAI	2.33	0.61
1:A:102:HIS:HB3	1:A:138:ARG:NH2	2.11	0.61
1:G:23:TRP:HA	1:G:30:ASN:HB3	1.83	0.60
1:H:86:THR:OG1	1:H:111:PHE:O	2.15	0.60
1:A:236:ASN:OD1	1:B:138:ARG:NH2	2.34	0.60
1:B:69:GLN:HB3	1:B:98:ASN:HB3	1.82	0.60
1:D:182:ASN:HA	1:D:243:SER:HA	1.84	0.60
1:G:80:VAL:HG12	1:G:87:LEU:HD11	1.83	0.60
1:C:164:LYS:HE2	1:C:166:GLN:O	2.02	0.60
1:B:134:ASP:CG	1:B:135:ALA:N	2.55	0.60
1:G:142:LEU:HD12	1:G:202:PHE:HB2	1.82	0.60
1:B:29:GLU:HA	1:B:34:ARG:HD3	1.82	0.60
1:D:189:SER:HA	1:D:197:GLU:HA	1.84	0.60
1:E:124:ARG:HG3	1:E:145:GLY:HA3	1.83	0.60
1:B:59:TYR:HE1	1:B:61:HIS:CE1	2.20	0.60
1:E:144:ILE:HB	1:E:204:ILE:HD12	1.84	0.60
1:F:124:ARG:CZ	1:F:207:PRO:HG3	2.31	0.60
1:G:105:VAL:O	1:G:132:HIS:HA	2.02	0.60
1:C:81:PHE:HA	1:C:162:ILE:HD13	1.84	0.59
1:G:116:GLU:HA	1:G:228:GLN:HG3	1.83	0.59
1:G:21:THR:HG22	1:G:22:LYS:H	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:GLU:HA	1:A:34:ARG:NH1	2.17	0.59
1:F:172:ALA:HB1	1:F:175:ALA:HB3	1.85	0.59
1:G:62:THR:OG1	1:H:61:HIS:O	2.06	0.59
1:D:88:LYS:NZ	3:D:303:7L3:OAM	2.35	0.59
1:G:130:PHE:HB2	1:G:142:LEU:HB3	1.83	0.59
1:B:179:LYS:HE2	1:B:247:ARG:HE	1.68	0.59
1:H:164:LYS:NZ	1:H:166:GLN:HB3	2.17	0.59
1:A:91:PHE:HE2	1:A:109[B]:VAL:HG12	1.67	0.59
1:F:111:PHE:HD2	1:F:126:LEU:HD11	1.68	0.59
1:C:86:THR:OG1	1:C:111:PHE:O	2.21	0.59
1:F:177:LEU:HB2	1:F:247:ARG:NH2	2.17	0.59
1:G:49:LYS:HB3	1:H:198:GLY:HA3	1.85	0.58
1:E:134:ASP:OD1	1:E:138:ARG:N	2.36	0.58
1:A:126:LEU:HB3	1:A:146:PHE:HB2	1.85	0.58
1:C:157:PRO:HB2	1:C:171:VAL:HG22	1.85	0.58
1:G:87:LEU:HD22	1:G:221:ARG:HD2	1.84	0.58
1:G:97:ILE:HD11	1:G:142:LEU:HD22	1.85	0.58
1:C:148:GLU:HA	1:C:210:VAL:O	2.03	0.58
1:D:64:ASP:OD1	1:D:183:TYR:OH	2.16	0.58
1:F:63:GLN:OE1	1:F:63:GLN:N	2.36	0.58
1:H:134:ASP:OD1	1:H:138:ARG:N	2.37	0.58
1:H:151:GLU:HG2	1:H:211:SER:CB	2.34	0.58
1:B:111:PHE:HB2	1:B:222:MET:CE	2.33	0.58
1:E:115:MET:HG3	1:E:127:SER:HB3	1.85	0.58
1:G:110:HIS:NE2	1:G:129:HIS:HB2	2.18	0.58
1:C:214:GLN:O	1:C:218:ILE:HG23	2.04	0.58
1:F:110:HIS:NE2	1:F:129:HIS:HB2	2.19	0.58
1:G:110:HIS:CE1	1:G:129:HIS:HB2	2.39	0.58
1:A:153:PRO:HA	1:A:156:ASP:OD2	2.04	0.57
1:B:149:GLY:N	1:B:210:VAL:O	2.31	0.57
1:A:104:TYR:HB3	1:A:132:HIS:HB3	1.87	0.57
1:F:95:ASN:HB2	1:F:106:LEU:HB3	1.86	0.57
1:A:165:LYS:HG3	1:A:166:GLN:H	1.68	0.57
1:G:150:LYS:O	1:G:209:GLU:HG2	2.04	0.57
1:H:23:TRP:HA	1:H:30:ASN:CB	2.34	0.57
1:G:60:TYR:HH	1:H:99:TYR:HH	1.49	0.57
1:A:173:LEU:HA	1:A:176:PHE:HB2	1.85	0.57
1:F:52:SER:O	1:F:229:ARG:NH1	2.38	0.57
1:B:166:GLN:HG3	1:B:167:ASN:H	1.70	0.57
1:H:183:TYR:CZ	1:H:242:SER:HB3	2.40	0.57
1:A:67:ASP:HB2	1:A:99:TYR:CE1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:116:GLU:HG2	1:D:228:GLN:HG3	1.87	0.57
1:C:133:LYS:HD3	3:C:303:7L3:CAV	2.35	0.56
1:C:32:PRO:HA	1:C:35:TRP:CD2	2.39	0.56
1:E:191:THR:O	1:E:229:ARG:HD2	2.05	0.56
1:B:133:LYS:HD2	3:B:302:7L3:CAW	2.35	0.56
1:C:58:HIS:CB	1:D:100:ARG:HH12	2.17	0.56
1:G:130:PHE:HB3	1:G:132:HIS:HE1	1.70	0.56
1:B:109:VAL:HA	1:B:129:HIS:O	2.05	0.56
1:A:62:THR:HG1	1:A:63:GLN:H	1.52	0.56
1:B:185:HIS:HE1	1:B:200:ALA:HB1	1.70	0.56
1:D:183:TYR:N	1:D:242:SER:O	2.33	0.56
1:A:159:LEU:HD23	1:A:217:GLU:OE1	2.05	0.56
1:H:80:VAL:HG12	1:H:162:ILE:HD13	1.87	0.56
1:A:172:ALA:HB1	1:A:175:ALA:HB3	1.87	0.56
1:B:76:LYS:HG3	1:B:170:GLU:HG2	1.86	0.56
1:C:218:ILE:O	1:C:222:MET:N	2.39	0.56
1:G:138:ARG:NH1	1:H:236:ASN:OD1	2.38	0.56
1:G:183:TYR:CZ	1:G:242:SER:HB3	2.41	0.56
1:A:219:LYS:HE3	1:A:224:ASN:HD22	1.71	0.56
1:D:115:MET:HG3	1:D:127:SER:HB3	1.86	0.56
1:G:235:TYR:HA	1:H:138:ARG:HH21	1.71	0.56
1:C:142:LEU:HD21	1:C:204:ILE:HD11	1.88	0.56
1:F:116:GLU:OE1	1:F:129:HIS:HE1	1.89	0.56
1:D:119:ILE:O	1:D:122:LYS:N	2.38	0.56
1:D:35:TRP:CH2	1:D:193:PRO:HD3	2.40	0.56
1:B:91:PHE:N	1:B:107:ASP:O	2.39	0.55
1:D:146:PHE:HA	1:D:208:LEU:O	2.06	0.55
3:H:302:7L3:SAA	3:H:302:7L3:OAO	2.64	0.55
1:A:132:HIS:HB2	1:A:140:LEU:HB3	1.88	0.55
1:B:217:GLU:O	1:B:221:ARG:HG2	2.07	0.55
1:C:110:HIS:CE1	1:C:129:HIS:HB2	2.41	0.55
1:F:32:PRO:HA	1:F:35:TRP:CD2	2.41	0.55
1:A:83:THR:HG21	1:A:88:LYS:HD3	1.88	0.55
1:C:236:ASN:O	1:D:187:ASN:ND2	2.30	0.55
1:D:179:LYS:NZ	1:D:245:GLU:OE2	2.39	0.55
1:H:164:LYS:HZ2	1:H:166:GLN:HB3	1.71	0.55
1:B:190:LEU:HD22	3:B:302:7L3:CAE	2.36	0.55
1:C:116:GLU:HG2	1:C:228:GLN:HG3	1.87	0.55
1:G:75:SER:O	1:G:171:VAL:N	2.25	0.55
1:C:97:ILE:O	1:C:104:TYR:N	2.24	0.55
1:F:116:GLU:HA	1:F:228:GLN:HG3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:ILE:HG23	1:D:239:ILE:HD13	1.87	0.55
1:E:214:GLN:O	1:E:218:ILE:HG13	2.07	0.55
1:H:219:LYS:HB3	1:H:224:ASN:HA	1.88	0.55
1:B:110:HIS:HE1	1:B:131:VAL:HG21	1.72	0.55
1:B:55:ASN:HB2	1:B:118:LEU:HD12	1.88	0.55
1:B:82:PHE:CE2	1:B:222:MET:HA	2.41	0.54
1:D:69:GLN:O	1:D:97:ILE:HG23	2.08	0.54
1:D:178:PRO:HD3	1:D:208:LEU:HD21	1.88	0.54
1:E:112:HIS:ND1	1:E:116:GLU:OE2	2.41	0.54
1:A:62:THR:OG1	1:A:63:GLN:N	2.40	0.54
1:B:71:LYS:HD2	1:B:96:HIS:HB2	1.89	0.54
1:C:71:LYS:HB2	1:C:96:HIS:HB2	1.88	0.54
1:F:24:ASP:OD1	1:F:30:ASN:HB2	2.07	0.54
1:E:130:PHE:O	1:E:141:VAL:HA	2.07	0.54
1:F:99:TYR:CZ	1:F:202:PHE:HZ	2.25	0.54
1:A:158:ILE:HD11	1:A:171:VAL:HG22	1.90	0.54
1:C:182:ASN:HB3	1:C:241:LYS:NZ	2.23	0.54
1:F:132:HIS:HB2	1:F:140:LEU:HB3	1.90	0.54
1:G:69:GLN:HG3	1:G:70:PHE:H	1.73	0.54
1:C:187:ASN:ND2	1:D:236:ASN:O	2.39	0.54
1:D:154:ASN:O	1:D:157:PRO:HD2	2.08	0.54
1:A:112:HIS:O	1:A:126:LEU:HD12	2.07	0.54
1:B:82:PHE:CD2	1:B:222:MET:HA	2.43	0.54
1:G:111:PHE:O	1:G:227:ASN:ND2	2.40	0.54
1:G:161:GLY:HA2	1:G:168:PHE:CE1	2.43	0.54
1:A:234:ASP:OD1	1:A:235:TYR:N	2.40	0.54
1:B:116:GLU:HA	1:B:228:GLN:HG3	1.89	0.54
1:C:164:LYS:NZ	1:C:167:ASN:O	2.23	0.54
1:C:67:ASP:OD1	1:D:60:TYR:OH	2.25	0.54
1:B:182:ASN:N	1:B:206:GLU:OE1	2.34	0.53
1:B:215:LEU:O	1:B:219:LYS:HG3	2.08	0.53
1:D:130:PHE:HB3	1:D:132:HIS:CE1	2.38	0.53
1:H:191:THR:N	3:H:302:7L3:OAH	2.33	0.53
1:A:179:LYS:HG3	1:A:247:ARG:HB2	1.89	0.53
3:A:302:7L3:OAO	3:A:302:7L3:SAA	2.66	0.53
1:G:106:LEU:HD11	1:G:108:ASN:O	2.08	0.53
1:G:142:LEU:CD1	1:G:202:PHE:HB2	2.37	0.53
1:H:162:ILE:HG21	1:H:221:ARG:HE	1.72	0.53
1:A:55:ASN:HA	1:A:118:LEU:HB2	1.90	0.53
1:D:63:GLN:N	1:D:63:GLN:OE1	2.41	0.53
1:F:79:ALA:O	1:F:89:ALA:HA	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:152:ASN:HB2	1:G:209:GLU:O	2.07	0.53
1:A:67:ASP:HB2	1:A:99:TYR:HE1	1.73	0.53
1:G:56:ILE:HG23	1:G:239:ILE:HD13	1.91	0.53
1:C:35:TRP:CH2	1:C:193:PRO:HD3	2.43	0.53
1:D:63:GLN:O	1:D:242:SER:OG	2.15	0.53
1:E:186:PHE:HB3	1:E:239:ILE:HG12	1.91	0.53
1:F:148:GLU:HA	1:F:210:VAL:O	2.08	0.53
3:D:303:7L3:SAA	3:D:303:7L3:OAO	2.67	0.53
1:E:151:GLU:O	1:E:153:PRO:HD3	2.09	0.53
1:E:35:TRP:CZ3	1:E:193:PRO:HD3	2.43	0.53
1:A:179:LYS:N	1:A:247:ARG:HH21	2.06	0.53
1:B:110:HIS:CE1	1:B:131:VAL:HG21	2.44	0.53
1:A:119:ILE:O	1:A:122:LYS:HB2	2.09	0.53
1:B:115:MET:HG3	1:B:127:SER:HB3	1.91	0.53
1:C:238:VAL:HG21	1:D:186:PHE:HA	1.90	0.52
1:F:181:ILE:HD13	1:F:246:THR:HG23	1.90	0.52
1:G:67:ASP:HB2	1:G:99:TYR:CD1	2.44	0.52
1:H:219:LYS:O	1:H:223:LYS:N	2.43	0.52
1:B:87:LEU:H	1:B:222:MET:CE	2.21	0.52
1:H:109:VAL:HA	1:H:129:HIS:O	2.09	0.52
1:H:105:VAL:O	1:H:132:HIS:HA	2.10	0.52
1:H:115:MET:HG3	1:H:127:SER:HB3	1.92	0.52
1:B:77:PRO:HD3	1:B:171:VAL:HG23	1.91	0.52
1:D:191:THR:OG1	3:D:303:7L3:NAI	2.42	0.52
1:G:35:TRP:CZ3	1:G:193:PRO:HD3	2.45	0.52
1:C:112:HIS:HA	1:C:227:ASN:HD21	1.74	0.52
1:F:164:LYS:HE2	1:F:166:GLN:OE1	2.10	0.52
1:C:110:HIS:HD2	1:C:112:HIS:NE2	2.07	0.52
1:D:67:ASP:N	1:D:67:ASP:OD1	2.40	0.52
3:G:303:7L3:SAA	3:G:303:7L3:OAO	2.68	0.52
1:G:68:LEU:HA	1:G:98:ASN:O	2.10	0.52
1:D:124:ARG:CZ	1:D:207:PRO:HB3	2.40	0.51
1:B:204:ILE:HG22	1:B:206:GLU:H	1.75	0.51
1:C:128:ALA:O	1:C:143:ALA:HA	2.11	0.51
1:C:76:LYS:HG3	1:C:170:GLU:HG2	1.91	0.51
1:A:150:LYS:H	1:A:209:GLU:HG2	1.75	0.51
1:B:159:LEU:HD21	1:B:218:ILE:HG12	1.92	0.51
1:D:51:GLN:HB3	1:D:232:GLN:HG3	1.93	0.51
1:F:147:GLU:HG3	1:F:209:GLU:HG3	1.92	0.51
1:B:83:THR:N	1:B:86:THR:O	2.30	0.51
1:H:157:PRO:HA	1:H:169:LYS:HE3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:216:ALA:HA	1:H:219:LYS:HZ1	1.75	0.51
1:B:88:LYS:HE3	1:B:108:ASN:CB	2.40	0.51
1:E:172:ALA:CB	1:E:175:ALA:HB3	2.40	0.51
1:F:146:PHE:HB3	1:F:210:VAL:HG12	1.93	0.51
1:E:51:GLN:OE1	1:E:229:ARG:NH2	2.38	0.51
1:B:88:LYS:HE3	1:B:108:ASN:CG	2.31	0.51
1:F:35:TRP:CZ3	1:F:193:PRO:HD3	2.46	0.51
1:E:26:LYS:C	1:E:28:LYS:H	2.13	0.51
1:F:105:VAL:O	1:F:132:HIS:HA	2.10	0.51
1:G:30:ASN:O	1:G:35:TRP:NE1	2.44	0.51
1:H:110:HIS:NE2	1:H:129:HIS:HB2	2.25	0.51
1:H:52:SER:HB3	1:H:117:PHE:CZ	2.45	0.51
1:A:164:LYS:HG2	1:A:165:LYS:O	2.11	0.51
1:B:83:THR:OG1	1:B:84:HIS:ND1	2.43	0.51
1:D:64:ASP:HB2	1:D:242:SER:OG	2.11	0.51
1:G:133:LYS:HG2	1:G:134:ASP:O	2.11	0.51
1:G:115:MET:O	1:G:228:GLN:NE2	2.44	0.51
1:D:96:HIS:HB3	1:D:103:ASP:OD1	2.10	0.50
1:A:119:ILE:HD12	1:A:124:ARG:HH21	1.76	0.50
1:B:102:HIS:HB3	1:B:134:ASP:CG	2.30	0.50
1:B:32:PRO:HA	1:B:35:TRP:CD2	2.46	0.50
1:C:63:GLN:OE1	1:G:61:HIS:ND1	2.44	0.50
1:F:217:GLU:OE2	1:F:221:ARG:NH1	2.44	0.50
1:A:152:ASN:HB2	1:A:209:GLU:O	2.11	0.50
1:D:183:TYR:CE1	1:D:242:SER:HB3	2.46	0.50
1:E:154:ASN:C	1:E:157:PRO:HD2	2.32	0.50
1:F:32:PRO:HA	1:F:35:TRP:CE2	2.47	0.50
3:F:302:7L3:OAO	3:F:302:7L3:SAA	2.70	0.50
1:G:104:TYR:HB3	1:G:132:HIS:HB3	1.92	0.50
1:C:58:HIS:HB2	1:D:100:ARG:HH12	1.76	0.50
1:E:146:PHE:HD2	1:E:210:VAL:HB	1.77	0.50
1:H:172:ALA:HB1	1:H:175:ALA:HB3	1.93	0.50
1:D:69:GLN:HB3	1:D:98:ASN:HB3	1.93	0.50
1:G:96:HIS:HB3	1:G:103:ASP:OD1	2.11	0.50
1:A:29:GLU:HA	1:A:34:ARG:HH12	1.75	0.49
1:H:78:LYS:HB2	1:H:90:SER:OG	2.11	0.49
1:A:147:GLU:HG3	1:A:148:GLU:HG2	1.94	0.49
1:D:179:LYS:O	1:D:246:THR:N	2.38	0.49
1:E:29:GLU:HA	1:E:34:ARG:NH2	2.28	0.49
1:E:138:ARG:HD2	1:F:235:TYR:HA	1.94	0.49
1:E:148:GLU:OE2	1:E:212:ALA:HA	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLN:HB3	1:A:232:GLN:HG3	1.93	0.49
1:C:116:GLU:HA	1:C:228:GLN:HG3	1.94	0.49
1:D:82:PHE:CE1	1:D:222:MET:HA	2.47	0.49
1:E:56:ILE:N	1:E:118:LEU:O	2.33	0.49
1:G:60:TYR:OH	1:H:99:TYR:OH	2.21	0.49
1:A:146:PHE:HA	1:A:208:LEU:O	2.13	0.49
1:G:30:ASN:HA	1:G:38:LEU:HD11	1.95	0.49
1:H:216:ALA:HA	1:H:219:LYS:NZ	2.27	0.49
1:H:104:TYR:HB3	1:H:132:HIS:HB3	1.94	0.49
1:C:183:TYR:CZ	1:C:242:SER:HB3	2.48	0.49
1:E:152:ASN:H	1:E:214:GLN:HE22	1.59	0.49
1:F:112:HIS:HA	1:F:227:ASN:ND2	2.28	0.49
1:A:130:PHE:HB3	1:A:132:HIS:CE1	2.48	0.49
1:B:76:LYS:HZ3	1:B:170:GLU:HG3	1.76	0.49
1:C:23:TRP:HB3	1:C:42:PHE:CE2	2.47	0.49
1:D:133:LYS:NZ	1:D:134:ASP:O	2.44	0.49
1:D:70:PHE:CE2	1:D:246:THR:HG23	2.48	0.49
1:D:179:LYS:HG3	1:D:247:ARG:HB2	1.95	0.49
1:D:30:ASN:HA	1:D:38:LEU:HD11	1.95	0.49
1:F:71:LYS:HB2	1:F:96:HIS:HB2	1.95	0.49
1:F:23:TRP:CZ2	1:F:84:HIS:HB3	2.48	0.49
1:B:104:TYR:HB3	1:B:132:HIS:HB3	1.95	0.49
1:E:77:PRO:HD2	1:E:168:PHE:HB3	1.95	0.49
1:F:125:PRO:HD2	1:F:146:PHE:O	2.13	0.49
1:G:106:LEU:HD12	1:G:131:VAL:O	2.13	0.49
1:H:23:TRP:HA	1:H:30:ASN:HB3	1.94	0.49
1:H:186:PHE:HD2	1:H:239:ILE:HG23	1.77	0.49
3:C:303:7L3:OAO	3:C:303:7L3:SAA	2.70	0.48
1:F:71:LYS:HD3	1:F:103:ASP:OD2	2.13	0.48
1:H:153:PRO:O	1:H:157:PRO:HD3	2.13	0.48
1:A:157:PRO:HB2	1:A:170:GLU:O	2.13	0.48
1:B:178:PRO:HD3	1:B:208:LEU:HD21	1.96	0.48
1:F:130:PHE:N	1:F:130:PHE:CD1	2.81	0.48
1:G:22:LYS:O	1:G:30:ASN:ND2	2.46	0.48
3:B:302:7L3:OAO	3:B:302:7L3:SAA	2.70	0.48
1:C:183:TYR:CE1	1:C:242:SER:HB3	2.47	0.48
1:D:116:GLU:HA	1:D:228:GLN:HG3	1.95	0.48
1:F:146:PHE:HA	1:F:208:LEU:O	2.13	0.48
1:G:54:ILE:HG22	1:G:232:GLN:O	2.13	0.48
1:H:36:ASP:OD1	1:H:36:ASP:N	2.43	0.48
1:A:79:ALA:HA	1:A:168:PHE:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:142:LEU:HD12	1:B:202:PHE:HB2	1.95	0.48
1:H:32:PRO:HA	1:H:35:TRP:CG	2.49	0.48
1:C:192:ALA:O	1:C:195:CYS:N	2.41	0.48
1:D:182:ASN:H	1:D:206:GLU:CD	2.17	0.48
1:E:33:HIS:C	1:E:46:LYS:HZ1	2.16	0.48
1:C:112:HIS:HA	1:C:227:ASN:ND2	2.29	0.48
1:E:223:LYS:N	1:E:223:LYS:HD2	2.29	0.48
1:E:78:LYS:HB2	1:E:90:SER:HB2	1.95	0.48
1:F:111:PHE:O	1:F:227:ASN:ND2	2.46	0.48
1:B:179:LYS:HG2	1:B:247:ARG:NE	2.29	0.48
1:E:23:TRP:CZ3	1:E:193:PRO:HD2	2.49	0.48
1:E:36:ASP:OD2	1:E:37:LYS:NZ	2.33	0.48
1:A:147:GLU:HG3	1:A:148:GLU:N	2.28	0.47
1:E:190:LEU:O	1:E:229:ARG:NH1	2.47	0.47
1:E:58:HIS:CD2	1:F:100:ARG:HH12	2.31	0.47
1:A:67:ASP:N	1:A:67:ASP:OD1	2.46	0.47
1:B:220:LYS:HA	1:B:224:ASN:H	1.79	0.47
1:C:102:HIS:HB2	1:C:104:TYR:CE1	2.49	0.47
1:C:35:TRP:CE3	1:C:193:PRO:HG3	2.48	0.47
1:D:28:LYS:O	1:D:33:HIS:CE1	2.67	0.47
1:F:147:GLU:O	1:F:209:GLU:HG3	2.15	0.47
1:H:23:TRP:O	1:H:85:HIS:NE2	2.35	0.47
1:A:156:ASP:N	1:A:157:PRO:HD2	2.29	0.47
1:A:164:LYS:HZ1	1:A:167:ASN:HB2	1.79	0.47
1:C:53:PRO:HA	1:C:117:PHE:CE1	2.49	0.47
1:D:154:ASN:HB2	1:D:176:PHE:CE1	2.49	0.47
1:E:118:LEU:CD2	1:E:123:THR:HG22	2.44	0.47
1:E:29:GLU:HA	1:E:34:ARG:HH22	1.79	0.47
1:F:42:PHE:CD1	1:F:42:PHE:N	2.81	0.47
1:F:110:HIS:CE1	3:F:302:7L3:SAA	3.07	0.47
1:F:67:ASP:OD1	1:F:68:LEU:N	2.45	0.47
1:A:139:LEU:HD21	3:A:302:7L3:CAV	2.45	0.47
1:D:112:HIS:O	1:D:126:LEU:HD12	2.14	0.47
1:G:76:LYS:HA	1:G:170:GLU:HA	1.96	0.47
1:G:115:MET:HG3	5:G:302:CL:CL	2.51	0.47
1:H:115:MET:HA	1:H:127:SER:HB3	1.96	0.47
1:B:166:GLN:HG3	1:B:167:ASN:N	2.29	0.47
1:D:59:TYR:HB3	1:D:239:ILE:HB	1.96	0.47
1:D:82:PHE:HZ	1:D:85:HIS:HA	1.80	0.47
3:E:302:7L3:SAA	3:E:302:7L3:OAO	2.72	0.47
1:E:82:PHE:C	1:E:82:PHE:CD1	2.88	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:95:ASN:CG	1:G:173:LEU:HD12	2.35	0.47
1:G:181:ILE:O	1:G:181:ILE:HG13	2.14	0.47
1:E:114:PRO:HD2	1:E:228:GLN:OE1	2.15	0.47
1:F:55:ASN:ND2	1:F:57:GLU:HG3	2.30	0.47
1:F:63:GLN:O	1:F:242:SER:OG	2.32	0.47
1:A:220:LYS:NZ	1:A:224:ASN:OD1	2.36	0.47
1:D:109:VAL:HA	1:D:129:HIS:O	2.15	0.47
1:D:184:TYR:HA	1:D:240:ILE:O	2.15	0.47
1:D:53:PRO:O	1:D:232:GLN:N	2.30	0.47
1:G:151:GLU:HG2	1:G:211:SER:HG	1.79	0.47
1:G:59:TYR:CE2	1:G:184:TYR:HE2	2.33	0.47
1:G:174:ASP:OD1	1:G:174:ASP:N	2.46	0.47
1:A:51:GLN:OE1	1:A:229:ARG:NH1	2.43	0.47
1:C:219:LYS:HB3	1:C:224:ASN:HA	1.97	0.47
1:F:178:PRO:HD3	1:F:208:LEU:HD21	1.97	0.47
1:B:168:PHE:CD1	1:B:168:PHE:N	2.84	0.46
1:B:140:LEU:HG	1:B:202:PHE:HE1	1.79	0.46
1:F:215:LEU:O	1:F:219:LYS:HG3	2.15	0.46
1:C:23:TRP:HB3	1:C:42:PHE:HE2	1.80	0.46
1:E:34:ARG:HA	1:E:36:ASP:OD1	2.15	0.46
1:C:102:HIS:HB2	1:C:104:TYR:HE1	1.79	0.46
1:B:42:PHE:N	1:B:42:PHE:CD1	2.81	0.46
1:C:62:THR:OG1	1:C:63:GLN:N	2.48	0.46
1:C:96:HIS:CE1	1:C:105:VAL:HG22	2.50	0.46
1:D:77:PRO:HD2	1:D:168:PHE:HB2	1.98	0.46
1:E:116:GLU:HG2	1:E:228:GLN:HG3	1.96	0.46
1:E:184:TYR:OH	1:E:205:GLU:OE2	2.28	0.46
1:H:77:PRO:HD2	1:H:168:PHE:HB2	1.97	0.46
1:B:210:VAL:HG23	1:B:214:GLN:OE1	2.15	0.46
1:D:56:ILE:N	1:D:118:LEU:O	2.42	0.46
1:A:112:HIS:O	1:A:127:SER:N	2.35	0.46
1:E:116:GLU:HA	1:E:228:GLN:HG3	1.97	0.46
1:E:72:TYR:OH	1:E:132:HIS:NE2	2.38	0.46
1:G:183:TYR:CE1	1:G:242:SER:HB3	2.51	0.46
1:G:127:SER:OG	5:G:302:CL:CL	2.58	0.46
1:D:114:PRO:HA	1:D:125:PRO:O	2.16	0.46
1:D:204:ILE:HG22	1:D:206:GLU:N	2.26	0.46
1:E:125:PRO:HD2	1:E:146:PHE:O	2.15	0.46
1:E:52:SER:O	1:E:229:ARG:NH1	2.49	0.46
1:G:238:VAL:HG21	1:H:186:PHE:HA	1.97	0.46
1:G:49:LYS:HB3	1:H:198:GLY:CA	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:215:LEU:O	1:H:219:LYS:NZ	2.48	0.46
1:A:63:GLN:HG2	1:A:64:ASP:N	2.30	0.46
1:C:235:TYR:CD2	1:D:138:ARG:HG2	2.51	0.46
1:F:54:ILE:HG12	1:F:117:PHE:HE1	1.79	0.46
1:G:89:ALA:O	1:G:108:ASN:HA	2.16	0.46
1:D:146:PHE:CE2	1:D:208:LEU:HB2	2.50	0.45
1:F:181:ILE:HG21	1:F:246:THR:HG21	1.98	0.45
1:G:151:GLU:HG2	1:G:211:SER:OG	2.16	0.45
1:G:56:ILE:HG23	1:G:239:ILE:HG21	1.99	0.45
1:D:151:GLU:OE2	1:D:213:LYS:NZ	2.39	0.45
1:D:44:VAL:HG13	1:D:195:CYS:HB2	1.98	0.45
1:D:42:PHE:N	1:D:42:PHE:CD1	2.83	0.45
1:G:74:ALA:HA	1:G:171:VAL:O	2.15	0.45
1:A:220:LYS:C	1:A:223:LYS:H	2.20	0.45
1:B:92:GLU:OE1	1:B:93:PRO:HD2	2.17	0.45
1:H:190:LEU:HA	1:H:190:LEU:HD23	1.70	0.45
1:H:125:PRO:HB2	1:H:215:LEU:HD11	1.98	0.45
1:A:112:HIS:CD2	1:A:129:HIS:CE1	3.05	0.45
1:C:29:GLU:CA	1:C:34:ARG:HE	2.28	0.45
1:D:117:PHE:HD2	5:D:302:CL:CL	2.37	0.45
1:C:86:THR:HG21	4:C:302:GOL:H12	1.99	0.45
1:D:182:ASN:N	1:D:206:GLU:OE1	2.50	0.45
1:B:151:GLU:OE1	1:B:151:GLU:N	2.49	0.45
1:D:25:TYR:HE1	1:D:229:ARG:HA	1.81	0.45
1:G:118:LEU:HG	1:G:123:THR:HG22	1.99	0.45
1:H:178:PRO:O	1:H:246:THR:OG1	2.18	0.45
1:B:35:TRP:CZ3	1:B:193:PRO:HD3	2.52	0.45
1:D:35:TRP:CZ2	1:D:193:PRO:HD3	2.52	0.45
1:E:82:PHE:CE2	1:E:223:LYS:HD3	2.52	0.45
1:B:150:LYS:O	1:B:211:SER:N	2.50	0.45
1:C:132:HIS:HB2	1:C:140:LEU:HB3	1.99	0.45
1:A:119:ILE:CD1	1:A:205:GLU:HA	2.42	0.45
1:A:26:LYS:HD3	1:A:26:LYS:HA	1.72	0.45
1:D:133:LYS:HA	1:D:138:ARG:O	2.17	0.45
1:F:177:LEU:C	1:F:247:ARG:HH22	2.20	0.45
1:B:112:HIS:O	1:B:126:LEU:HD12	2.17	0.44
1:D:130:PHE:HB2	1:D:142:LEU:HB3	1.99	0.44
1:D:150:LYS:O	1:D:211:SER:N	2.49	0.44
1:D:181:ILE:HG13	1:D:181:ILE:O	2.16	0.44
1:B:39:HIS:ND1	1:D:41:ASP:OD2	2.50	0.44
1:F:83:THR:O	1:F:86:THR:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:211:SER:OG	1:G:214:GLN:N	2.47	0.44
1:C:217:GLU:O	1:C:220:LYS:HB3	2.17	0.44
1:D:110:HIS:CE1	1:D:129:HIS:HB2	2.52	0.44
1:D:23:TRP:CD1	1:D:42:PHE:HE2	2.35	0.44
1:F:115:MET:HG3	1:F:127:SER:HB3	1.99	0.44
1:F:70:PHE:HB2	1:F:247:ARG:HG3	2.00	0.44
1:A:32:PRO:HA	1:A:35:TRP:CE2	2.52	0.44
1:B:227:ASN:OD1	1:B:227:ASN:N	2.51	0.44
1:A:76:LYS:HA	1:A:77:PRO:HD3	1.85	0.44
1:E:80:VAL:HB	1:E:161:GLY:O	2.17	0.44
1:F:88:LYS:HA	1:F:109:VAL:O	2.17	0.44
1:H:64:ASP:N	1:H:64:ASP:OD1	2.50	0.44
1:C:57:GLU:HG2	1:C:58:HIS:CE1	2.52	0.44
1:D:110:HIS:NE2	1:D:129:HIS:HB2	2.32	0.44
1:H:26:LYS:C	1:H:28:LYS:H	2.19	0.44
1:A:165:LYS:HA	1:A:168:PHE:CE2	2.53	0.44
1:A:78:LYS:HB3	1:A:78:LYS:HE2	1.80	0.44
1:B:184:TYR:O	1:B:202:PHE:HA	2.17	0.44
1:B:87:LEU:HD12	1:B:88:LYS:H	1.83	0.44
1:F:174:ASP:OD1	1:F:174:ASP:N	2.49	0.44
1:G:166:GLN:HB3	1:G:167:ASN:H	1.64	0.44
1:H:79:ALA:HB2	1:H:165:LYS:NZ	2.33	0.44
1:H:77:PRO:HA	1:H:90:SER:O	2.18	0.44
1:A:105:VAL:O	1:A:132:HIS:HA	2.18	0.44
1:B:106:LEU:HD12	1:B:131:VAL:O	2.17	0.44
1:B:105:VAL:O	1:B:132:HIS:HA	2.17	0.44
1:C:27:ASN:OD1	1:C:27:ASN:N	2.41	0.44
1:C:42:PHE:N	1:C:42:PHE:CD1	2.84	0.44
1:C:82:PHE:C	1:C:82:PHE:CD1	2.91	0.44
1:D:146:PHE:HB3	1:D:210:VAL:HG12	2.00	0.44
1:F:211:SER:HB2	1:F:213:LYS:NZ	2.33	0.44
1:C:184:TYR:HE1	1:C:205:GLU:HB2	1.82	0.44
1:D:82:PHE:CZ	1:D:85:HIS:HA	2.53	0.44
1:F:118:LEU:HG	1:F:123:THR:CG2	2.48	0.44
1:F:158:ILE:HD11	1:F:171:VAL:HG21	2.00	0.44
1:G:106:LEU:HA	1:G:131:VAL:O	2.18	0.44
1:G:139:LEU:HD13	1:G:190:LEU:HD21	2.00	0.44
1:B:51:GLN:OE1	1:B:195:CYS:HB3	2.17	0.44
1:D:177:LEU:CG	1:D:246:THR:HG21	2.48	0.44
1:D:116:GLU:HB2	5:D:302:CL:CL	2.54	0.44
1:G:204:ILE:HG22	1:G:206:GLU:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:TRP:HZ2	1:C:84:HIS:HB3	1.82	0.43
1:D:59:TYR:CE2	1:D:184:TYR:HE2	2.36	0.43
1:D:44:VAL:HA	1:D:47:SER:OG	2.17	0.43
1:G:105:VAL:HG12	1:G:133:LYS:O	2.18	0.43
1:A:130:PHE:HB3	1:A:132:HIS:HE1	1.82	0.43
1:A:219:LYS:HE3	1:A:224:ASN:ND2	2.33	0.43
1:A:22:LYS:NZ	1:A:30:ASN:OD1	2.43	0.43
1:D:86:THR:HA	1:D:222:MET:SD	2.57	0.43
1:E:118:LEU:HD13	1:E:121:ASN:HA	1.99	0.43
1:E:147:GLU:O	1:E:209:GLU:HA	2.17	0.43
1:F:181:ILE:CD1	1:F:244:ALA:H	2.30	0.43
1:H:27:ASN:OD1	1:H:27:ASN:N	2.50	0.43
1:B:119:ILE:O	1:B:122:LYS:N	2.38	0.43
1:C:104:TYR:HB3	1:C:132:HIS:CB	2.45	0.43
1:C:189:SER:HA	1:C:197:GLU:HA	2.00	0.43
1:C:34:ARG:O	1:C:38:LEU:HG	2.17	0.43
1:D:104:TYR:CD2	1:D:134:ASP:HB3	2.52	0.43
1:G:184:TYR:CE2	1:G:241:LYS:HD2	2.54	0.43
1:H:64:ASP:O	1:H:66:ALA:N	2.51	0.43
1:D:227:ASN:OD1	1:D:227:ASN:N	2.43	0.43
1:A:113:ALA:HB2	1:A:126:LEU:HD13	1.99	0.43
1:A:64:ASP:OD1	1:A:242:SER:OG	2.36	0.43
1:D:133:LYS:HG2	1:D:134:ASP:O	2.18	0.43
1:F:66:ALA:HA	1:F:244:ALA:HB1	2.01	0.43
1:G:113:ALA:HA	1:G:114:PRO:HA	1.87	0.43
1:A:35:TRP:CH2	1:A:193:PRO:HD3	2.53	0.43
1:C:100:ARG:HD2	1:D:58:HIS:CE1	2.53	0.43
1:C:234:ASP:O	1:D:138:ARG:NE	2.52	0.43
1:E:42:PHE:N	1:E:42:PHE:CD1	2.87	0.43
1:G:30:ASN:O	1:G:35:TRP:CD1	2.72	0.43
1:H:182:ASN:OD1	1:H:243:SER:HB3	2.19	0.43
1:E:31:GLY:HA3	1:E:33:HIS:CE1	2.54	0.43
1:F:192:ALA:HB3	3:F:302:7L3:NAD	2.34	0.43
1:A:189:SER:HA	1:A:197:GLU:HA	2.00	0.43
1:B:112:HIS:CD2	1:B:129:HIS:CE1	3.06	0.43
1:B:32:PRO:HA	1:B:35:TRP:CE2	2.54	0.43
1:C:105:VAL:O	1:C:132:HIS:HA	2.18	0.43
1:E:152:ASN:N	1:E:214:GLN:HE22	2.17	0.43
1:F:134:ASP:OD1	1:F:138:ARG:N	2.52	0.43
1:B:168:PHE:H	1:B:168:PHE:HD1	1.66	0.43
1:C:56:ILE:O	1:C:120:ASN:N	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:146:PHE:CD2	1:E:210:VAL:HB	2.54	0.43
1:F:114:PRO:O	1:F:228:GLN:HB2	2.19	0.43
1:H:59:TYR:CD1	1:H:59:TYR:N	2.87	0.43
1:E:227:ASN:OD1	1:E:227:ASN:N	2.48	0.42
1:E:30:ASN:HA	1:E:38:LEU:HD11	2.00	0.42
1:F:105:VAL:HG12	1:F:133:LYS:O	2.19	0.42
1:F:54:ILE:O	1:F:117:PHE:HD1	2.02	0.42
1:F:62:THR:OG1	1:F:63:GLN:N	2.51	0.42
1:H:186:PHE:HB3	1:H:239:ILE:HG12	2.01	0.42
1:A:138:ARG:HG2	1:B:235:TYR:HD2	1.85	0.42
1:A:102:HIS:CG	1:A:138:ARG:HH12	2.37	0.42
1:B:225:SER:HA	1:B:226:PRO:HD3	1.77	0.42
1:D:163:GLN:O	1:D:164:LYS:HD2	2.19	0.42
1:A:150:LYS:HG2	1:A:151:GLU:N	2.33	0.42
1:A:96:HIS:HB3	1:A:103:ASP:OD1	2.20	0.42
1:G:219:LYS:HG2	1:G:222:MET:HE2	2.01	0.42
1:H:129:HIS:HA	1:H:142:LEU:O	2.20	0.42
1:H:53:PRO:HA	1:H:117:PHE:CE1	2.54	0.42
1:A:217:GLU:O	1:A:221:ARG:HG2	2.19	0.42
1:C:99:TYR:CZ	1:C:100:ARG:HG3	2.53	0.42
1:F:194:PRO:HG2	1:F:196:THR:OG1	2.19	0.42
1:G:115:MET:HG2	1:G:117:PHE:O	2.19	0.42
1:C:115:MET:HB2	1:C:124:ARG:HB2	2.01	0.42
1:G:104:TYR:CB	1:G:132:HIS:HB3	2.50	0.42
1:G:183:TYR:O	1:G:241:LYS:HA	2.20	0.42
1:B:151:GLU:O	1:B:153:PRO:HD3	2.20	0.42
1:B:112:HIS:HA	1:B:227:ASN:ND2	2.35	0.42
1:C:128:ALA:HB2	1:C:146:PHE:HE1	1.84	0.42
1:C:67:ASP:O	1:C:99:TYR:HD1	2.02	0.42
1:D:105:VAL:O	1:D:132:HIS:HA	2.20	0.42
1:D:81:PHE:N	1:D:81:PHE:CD1	2.87	0.42
1:E:161:GLY:HA2	1:E:168:PHE:CD1	2.54	0.42
1:E:225:SER:HA	1:E:226:PRO:HD3	1.74	0.42
1:G:110:HIS:HE1	1:G:131:VAL:CG2	2.32	0.42
1:H:230:PRO:O	1:H:232:GLN:NE2	2.42	0.42
1:A:117:PHE:CE1	1:A:186:PHE:HZ	2.37	0.42
1:C:32:PRO:HA	1:C:35:TRP:CD1	2.53	0.42
1:E:158:ILE:HD11	1:E:171:VAL:HG21	2.02	0.42
1:F:174:ASP:HA	1:F:247:ARG:HH11	1.84	0.42
1:F:55:ASN:HD21	1:F:57:GLU:HG3	1.85	0.42
1:H:35:TRP:CZ3	1:H:193:PRO:HD3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:HIS:CE1	1:A:129:HIS:HB2	2.55	0.42
1:A:51:GLN:OE1	1:A:229:ARG:NH2	2.52	0.42
1:C:108:ASN:HD21	1:C:131:VAL:HG21	1.85	0.42
1:C:32:PRO:HG3	1:C:35:TRP:CZ2	2.55	0.42
1:E:116:GLU:OE1	1:E:129:HIS:HE1	2.02	0.42
1:G:92:GLU:O	1:G:94:THR:N	2.50	0.42
1:H:165:LYS:O	1:H:165:LYS:HG2	2.20	0.42
1:A:66:ALA:O	1:A:69:GLN:NE2	2.52	0.42
1:C:73:ALA:O	1:C:173:LEU:HB2	2.20	0.42
1:C:190:LEU:HD23	1:C:190:LEU:HA	1.82	0.42
1:C:58:HIS:HB2	1:D:100:ARG:NH1	2.35	0.42
1:D:164:LYS:HG2	1:D:166:GLN:NE2	2.34	0.42
1:D:234:ASP:OD1	1:D:237:THR:N	2.42	0.42
1:E:191:THR:HA	1:E:229:ARG:HH11	1.83	0.42
1:F:129:HIS:CD2	1:F:143:ALA:HB2	2.55	0.42
1:B:181:ILE:HA	1:B:206:GLU:OE1	2.21	0.41
1:D:54:ILE:HG22	1:D:232:GLN:O	2.19	0.41
1:H:185:HIS:CE1	1:H:202:PHE:CZ	3.08	0.41
1:F:57:GLU:HG2	1:F:120:ASN:OD1	2.21	0.41
1:G:130:PHE:HB3	1:G:132:HIS:CE1	2.53	0.41
1:A:38:LEU:HA	1:A:38:LEU:HD23	1.88	0.41
1:B:55:ASN:HA	1:B:118:LEU:HB2	2.02	0.41
1:C:165:LYS:HD3	1:C:165:LYS:HA	1.93	0.41
1:D:73:ALA:O	1:D:173:LEU:HD12	2.20	0.41
1:D:185:HIS:HB3	1:D:240:ILE:HG13	2.02	0.41
1:G:87:LEU:HD12	1:G:88:LYS:H	1.85	0.41
1:A:89:ALA:O	1:A:108:ASN:HB2	2.20	0.41
1:C:182:ASN:HB3	1:C:241:LYS:HZ3	1.85	0.41
1:E:197:GLU:HB2	1:F:49:LYS:HG3	2.02	0.41
1:B:179:LYS:HG2	1:B:247:ARG:CZ	2.50	0.41
1:C:109[B]:VAL:HA	1:C:129:HIS:O	2.21	0.41
1:E:152:ASN:H	1:E:214:GLN:NE2	2.18	0.41
1:F:32:PRO:HA	1:F:35:TRP:CG	2.55	0.41
1:G:154:ASN:O	1:G:157:PRO:HD2	2.21	0.41
1:G:219:LYS:HE3	1:G:219:LYS:HB3	1.82	0.41
1:G:174:ASP:O	1:G:247:ARG:NH1	2.54	0.41
1:A:91:PHE:CE2	1:A:109[B]:VAL:HG12	2.52	0.41
1:A:234:ASP:OD1	1:A:237:THR:N	2.47	0.41
1:C:115:MET:HG3	1:C:127:SER:CB	2.46	0.41
1:E:129:HIS:HB3	1:E:141:VAL:HG11	2.02	0.41
1:E:204:ILE:HG22	1:E:206:GLU:N	2.27	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:79:ALA:HB2	1:H:165:LYS:HZ3	1.85	0.41
1:A:228:GLN:HG2	1:A:229:ARG:O	2.20	0.41
1:D:87:LEU:HD12	1:D:88:LYS:H	1.86	0.41
1:E:112:HIS:ND1	1:E:116:GLU:CD	2.74	0.41
1:E:138:ARG:HG2	1:F:235:TYR:CD2	2.56	0.41
1:F:149:GLY:HA3	1:F:209:GLU:OE2	2.21	0.41
1:D:217:GLU:O	1:D:221:ARG:HG2	2.21	0.41
1:F:111:PHE:HB2	1:F:222:MET:SD	2.61	0.41
1:G:75:SER:N	1:G:171:VAL:O	2.45	0.41
1:H:225:SER:HA	1:H:226:PRO:HD3	1.68	0.41
1:H:31:GLY:HA2	1:H:32:PRO:HD3	1.87	0.41
1:A:31:GLY:O	1:A:34:ARG:HB2	2.21	0.41
1:A:52:SER:HB3	1:A:117:PHE:CZ	2.56	0.41
1:A:106:LEU:HD12	1:A:131:VAL:O	2.21	0.41
1:B:83:THR:O	1:B:84:HIS:C	2.59	0.41
1:B:71:LYS:N	1:B:96:HIS:O	2.27	0.41
1:C:115:MET:HB3	1:C:124:ARG:H	1.85	0.41
1:C:185:HIS:HA	1:C:201:TRP:O	2.21	0.41
1:E:115:MET:CG	1:E:127:SER:HB3	2.51	0.41
1:F:104:TYR:HA	1:F:134:ASP:HA	2.02	0.41
1:G:21:THR:HG22	1:G:22:LYS:N	2.34	0.41
1:H:111:PHE:HA	1:H:128:ALA:HA	2.03	0.41
1:B:81:PHE:O	1:B:88:LYS:N	2.52	0.40
1:D:191:THR:O	1:D:229:ARG:HB2	2.21	0.40
1:G:191:THR:OG1	3:G:303:7L3:NAI	2.54	0.40
1:C:71:LYS:HD2	1:C:96:HIS:CB	2.45	0.40
1:F:49:LYS:HA	1:F:49:LYS:HD3	1.80	0.40
1:G:70:PHE:CD1	1:G:246:THR:HA	2.56	0.40
1:A:184:TYR:CE2	1:A:241:LYS:HD2	2.57	0.40
1:E:22:LYS:O	1:E:30:ASN:ND2	2.54	0.40
1:E:64:ASP:OD1	1:E:244:ALA:HA	2.21	0.40
1:F:159:LEU:O	1:F:162:ILE:HG22	2.22	0.40
1:G:174:ASP:O	1:G:247:ARG:HD3	2.22	0.40
1:B:36:ASP:HA	1:B:42:PHE:O	2.22	0.40
1:D:136:LYS:HD2	1:D:138:ARG:HH11	1.87	0.40
1:D:56:ILE:HG21	1:D:184:TYR:CD2	2.56	0.40
1:E:162:ILE:HD12	1:E:162:ILE:HA	1.90	0.40
1:F:181:ILE:HD11	1:F:244:ALA:H	1.87	0.40
1:G:32:PRO:HA	1:G:35:TRP:CE2	2.55	0.40
1:H:145:GLY:O	1:H:207:PRO:HA	2.21	0.40
1:C:215:LEU:O	1:C:219:LYS:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:PRO:HG3	1:D:91:PHE:CZ	2.56	0.40
1:E:75:SER:HB2	1:E:91:PHE:CE1	2.56	0.40
1:E:95:ASN:HB2	1:E:106:LEU:HB3	2.03	0.40
1:F:133:LYS:HD2	1:F:137:GLY:HA2	2.02	0.40
1:G:186:PHE:CE1	1:G:201:TRP:HB2	2.57	0.40
1:H:209:GLU:H	1:H:209:GLU:HG3	1.33	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:HIS:ND1	1:B:148:GLU:OE1[1_655]	2.14	0.06

## 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	226/234 (97%)	211 (93%)	15 (7%)	0	100	100
1	B	213/234 (91%)	200 (94%)	13 (6%)	0	100	100
1	C	225/234 (96%)	209 (93%)	16 (7%)	0	100	100
1	D	212/234 (91%)	198 (93%)	14 (7%)	0	100	100
1	E	225/234 (96%)	209 (93%)	16 (7%)	0	100	100
1	F	216/234 (92%)	202 (94%)	14 (6%)	0	100	100
1	G	217/234 (93%)	198 (91%)	19 (9%)	0	100	100
1	H	223/234 (95%)	208 (93%)	15 (7%)	0	100	100
All	All	1757/1872 (94%)	1635 (93%)	122 (7%)	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	203/208 (98%)	200 (98%)	3 (2%)	65	87
1	B	195/208 (94%)	187 (96%)	8 (4%)	30	64
1	C	202/208 (97%)	199 (98%)	3 (2%)	65	87
1	D	196/208 (94%)	191 (97%)	5 (3%)	46	77
1	E	202/208 (97%)	200 (99%)	2 (1%)	76	92
1	F	197/208 (95%)	193 (98%)	4 (2%)	55	82
1	G	199/208 (96%)	195 (98%)	4 (2%)	55	82
1	H	200/208 (96%)	196 (98%)	4 (2%)	55	82
All	All	1594/1664 (96%)	1561 (98%)	33 (2%)	55	81

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50[A]	SER
1	A	50[B]	SER
1	A	111	PHE
1	B	39	HIS
1	B	86	THR
1	B	109	VAL
1	B	138	ARG
1	B	168	PHE
1	B	218	ILE
1	B	238	VAL
1	B	243	SER
1	C	39	HIS
1	C	136	LYS
1	C	151	GLU
1	D	50	SER
1	D	81	PHE
1	D	131	VAL
1	D	176	PHE
1	D	184	TYR

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Mol	Chain	Res	Type
1	E	82	PHE
1	E	170	GLU
1	F	27	ASN
1	F	103	ASP
1	F	148	GLU
1	F	152	ASN
1	G	67	ASP
1	G	86	THR
1	G	136	LYS
1	G	220	LYS
1	H	59	TYR
1	H	63	GLN
1	H	209	GLU
1	H	243	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	108	ASN
1	E	95	ASN
1	E	96	HIS
1	F	55	ASN
1	G	95	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 10 are monoatomic - leaving 9 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$
3	7L3	H	302	2	19,25,25	3.31	7 (36%)	29,40,40	1.65	2 (6%)
3	7L3	D	303	2	19,25,25	3.30	7 (36%)	29,40,40	1.64	2 (6%)
3	7L3	C	303	2	19,25,25	3.30	7 (36%)	29,40,40	1.64	2 (6%)
3	7L3	G	303	2	19,25,25	3.30	7 (36%)	29,40,40	1.65	2 (6%)
4	GOL	C	302	-	5,5,5	0.25	0	5,5,5	0.28	0
3	7L3	B	302	2	19,25,25	3.30	7 (36%)	29,40,40	1.64	2 (6%)
3	7L3	A	302	2	19,25,25	3.52	8 (42%)	29,40,40	1.67	2 (6%)
3	7L3	F	302	2	19,25,25	3.50	8 (42%)	29,40,40	1.67	2 (6%)
3	7L3	E	302	2	19,25,25	3.29	7 (36%)	29,40,40	1.65	2 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	7L3	H	302	2	-	11/15/23/23	0/2/2/2
3	7L3	D	303	2	-	11/15/23/23	0/2/2/2
3	7L3	C	303	2	-	9/15/23/23	0/2/2/2
3	7L3	G	303	2	-	11/15/23/23	0/2/2/2
4	GOL	C	302	-	-	2/4/4/4	-
3	7L3	B	302	2	-	9/15/23/23	0/2/2/2
3	7L3	A	302	2	-	4/15/23/23	0/2/2/2
3	7L3	F	302	2	-	9/15/23/23	0/2/2/2
3	7L3	E	302	2	-	11/15/23/23	0/2/2/2

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	7L3	CAB-NAK	8.53	1.44	1.32
3	F	302	7L3	CAB-NAK	8.48	1.44	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	303	7L3	CAB-NAK	8.07	1.44	1.32
3	C	303	7L3	CAB-NAK	8.07	1.44	1.32
3	D	303	7L3	CAB-NAK	8.05	1.44	1.32
3	B	302	7L3	CAB-NAK	8.04	1.44	1.32
3	H	302	7L3	CAB-NAK	8.03	1.44	1.32
3	E	302	7L3	CAB-NAK	8.02	1.44	1.32
3	B	302	7L3	OAO-SAL	7.03	1.52	1.44
3	H	302	7L3	OAO-SAL	7.02	1.52	1.44
3	C	303	7L3	OAO-SAL	6.99	1.52	1.44
3	E	302	7L3	OAO-SAL	6.99	1.52	1.44
3	D	303	7L3	OAO-SAL	6.98	1.52	1.44
3	G	303	7L3	OAO-SAL	6.95	1.52	1.44
3	A	302	7L3	OAO-SAL	6.22	1.51	1.44
3	F	302	7L3	OAO-SAL	6.19	1.51	1.44
3	H	302	7L3	OAM-SAL	5.38	1.50	1.44
3	B	302	7L3	OAM-SAL	5.33	1.50	1.44
3	C	303	7L3	OAM-SAL	5.32	1.50	1.44
3	D	303	7L3	OAM-SAL	5.31	1.50	1.44
3	G	303	7L3	OAM-SAL	5.30	1.50	1.44
3	E	302	7L3	OAM-SAL	5.30	1.50	1.44
3	F	302	7L3	OAG-SAF	5.13	1.53	1.43
3	A	302	7L3	OAG-SAF	5.12	1.53	1.43
3	A	302	7L3	OAM-SAL	5.09	1.50	1.44
3	F	302	7L3	OAM-SAL	4.98	1.50	1.44
3	F	302	7L3	CAN-SAL	4.93	1.83	1.76
3	A	302	7L3	CAN-SAL	4.89	1.83	1.76
3	G	303	7L3	OAG-SAF	4.75	1.52	1.43
3	H	302	7L3	OAG-SAF	4.72	1.52	1.43
3	C	303	7L3	OAG-SAF	4.71	1.52	1.43
3	E	302	7L3	OAG-SAF	4.68	1.52	1.43
3	B	302	7L3	OAG-SAF	4.66	1.52	1.43
3	D	303	7L3	OAG-SAF	4.65	1.52	1.43
3	A	302	7L3	SAF-NAI	4.21	1.68	1.60
3	F	302	7L3	SAF-NAI	4.15	1.68	1.60
3	H	302	7L3	OAH-SAF	3.95	1.51	1.43
3	D	303	7L3	OAH-SAF	3.91	1.51	1.43
3	B	302	7L3	OAH-SAF	3.90	1.51	1.43
3	C	303	7L3	OAH-SAF	3.89	1.51	1.43
3	E	302	7L3	OAH-SAF	3.89	1.51	1.43
3	G	303	7L3	OAH-SAF	3.86	1.50	1.43
3	E	302	7L3	CAN-SAL	3.28	1.81	1.76
3	H	302	7L3	CAN-SAL	3.26	1.81	1.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	303	7L3	CAN-SAL	3.25	1.81	1.76
3	B	302	7L3	CAN-SAL	3.23	1.81	1.76
3	D	303	7L3	CAN-SAL	3.22	1.81	1.76
3	G	303	7L3	CAN-SAL	3.20	1.81	1.76
3	A	302	7L3	OAH-SAF	2.96	1.49	1.43
3	F	302	7L3	OAH-SAF	2.93	1.49	1.43
3	A	302	7L3	CAP-CAN	2.37	1.42	1.38
3	E	302	7L3	SAF-NAI	2.34	1.65	1.60
3	B	302	7L3	SAF-NAI	2.34	1.65	1.60
3	F	302	7L3	CAP-CAN	2.32	1.42	1.38
3	D	303	7L3	SAF-NAI	2.31	1.65	1.60
3	G	303	7L3	SAF-NAI	2.29	1.64	1.60
3	H	302	7L3	SAF-NAI	2.28	1.64	1.60
3	C	303	7L3	SAF-NAI	2.26	1.64	1.60

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	302	7L3	OAH-SAF-OAG	-5.71	109.36	118.76
3	A	302	7L3	OAH-SAF-OAG	-5.71	109.37	118.76
3	G	303	7L3	OAH-SAF-OAG	-5.68	109.42	118.76
3	H	302	7L3	OAH-SAF-OAG	-5.68	109.43	118.76
3	C	303	7L3	OAH-SAF-OAG	-5.68	109.43	118.76
3	D	303	7L3	OAH-SAF-OAG	-5.66	109.46	118.76
3	E	302	7L3	OAH-SAF-OAG	-5.65	109.47	118.76
3	B	302	7L3	OAH-SAF-OAG	-5.65	109.47	118.76
3	E	302	7L3	OAO-SAL-OAM	-5.40	109.45	117.21
3	B	302	7L3	OAO-SAL-OAM	-5.39	109.47	117.21
3	G	303	7L3	OAO-SAL-OAM	-5.38	109.48	117.21
3	H	302	7L3	OAO-SAL-OAM	-5.38	109.48	117.21
3	C	303	7L3	OAO-SAL-OAM	-5.37	109.50	117.21
3	D	303	7L3	OAO-SAL-OAM	-5.36	109.50	117.21
3	A	302	7L3	OAO-SAL-OAM	-5.32	109.56	117.21
3	F	302	7L3	OAO-SAL-OAM	-5.28	109.62	117.21

There are no chirality outliers.

All (77) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	302	GOL	C1-C2-C3-O3
3	B	302	7L3	CAP-CAN-SAL-NAK
3	A	302	7L3	CAP-CAN-SAL-NAK

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Mol	Chain	Res	Type	Atoms
3	F	302	7L3	CAP-CAN-SAL-NAK
3	B	302	7L3	CAT-CAN-SAL-NAK
3	A	302	7L3	CAT-CAN-SAL-NAK
3	C	303	7L3	CAP-CAN-SAL-NAK
3	F	302	7L3	CAT-CAN-SAL-NAK
4	C	302	GOL	O2-C2-C3-O3
3	E	302	7L3	CAP-CAN-SAL-NAK
3	C	303	7L3	CAT-CAN-SAL-NAK
3	E	302	7L3	CAT-CAN-SAL-NAK
3	H	302	7L3	CAB-NAK-SAL-OAM
3	B	302	7L3	CAB-NAK-SAL-OAM
3	G	303	7L3	CAP-CAN-SAL-NAK
3	D	303	7L3	CAP-CAN-SAL-NAK
3	G	303	7L3	CAB-NAK-SAL-OAM
3	D	303	7L3	CAB-NAK-SAL-OAM
3	D	303	7L3	CAT-CAN-SAL-NAK
3	G	303	7L3	CAT-CAN-SAL-NAK
3	E	302	7L3	CAB-NAK-SAL-OAM
3	A	302	7L3	CAB-NAK-SAL-OAM
3	H	302	7L3	CAT-CAN-SAL-OAO
3	H	302	7L3	CAP-CAN-SAL-NAK
3	H	302	7L3	CAP-CAN-SAL-OAO
3	G	303	7L3	CAT-CAN-SAL-OAO
3	D	303	7L3	CAT-CAN-SAL-OAO
3	D	303	7L3	CAP-CAN-SAL-OAO
3	G	303	7L3	CAP-CAN-SAL-OAO
3	D	303	7L3	CAQ-CAR-CAU-CAX
3	G	303	7L3	CAS-CAR-CAU-CAX
3	E	302	7L3	CAS-CAR-CAU-CAX
3	H	302	7L3	CAS-CAR-CAU-CAW
3	H	302	7L3	CAQ-CAR-CAU-CAW
3	C	303	7L3	CAQ-CAR-CAU-CAV
3	B	302	7L3	CAQ-CAR-CAU-CAX
3	D	303	7L3	CAS-CAR-CAU-CAX
3	C	303	7L3	CAS-CAR-CAU-CAV
3	B	302	7L3	CAS-CAR-CAU-CAX
3	E	302	7L3	CAQ-CAR-CAU-CAX
3	F	302	7L3	CAQ-CAR-CAU-CAW
3	G	303	7L3	CAQ-CAR-CAU-CAX
3	F	302	7L3	CAS-CAR-CAU-CAW
3	E	302	7L3	CAT-CAN-SAL-OAM
3	H	302	7L3	CAT-CAN-SAL-NAK

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Mol	Chain	Res	Type	Atoms
3	H	302	7L3	CAQ-CAR-CAU-CAV
3	H	302	7L3	CAS-CAR-CAU-CAV
3	H	302	7L3	CAS-CAR-CAU-CAX
3	D	303	7L3	CAQ-CAR-CAU-CAV
3	D	303	7L3	CAS-CAR-CAU-CAW
3	C	303	7L3	CAS-CAR-CAU-CAW
3	C	303	7L3	CAQ-CAR-CAU-CAW
3	C	303	7L3	CAQ-CAR-CAU-CAX
3	B	302	7L3	CAQ-CAR-CAU-CAV
3	B	302	7L3	CAS-CAR-CAU-CAW
3	B	302	7L3	CAQ-CAR-CAU-CAW
3	E	302	7L3	CAS-CAR-CAU-CAW
3	D	303	7L3	CAQ-CAR-CAU-CAW
3	G	303	7L3	CAS-CAR-CAU-CAV
3	G	303	7L3	CAS-CAR-CAU-CAW
3	E	302	7L3	CAS-CAR-CAU-CAV
3	H	302	7L3	CAQ-CAR-CAU-CAX
3	C	303	7L3	CAS-CAR-CAU-CAX
3	B	302	7L3	CAS-CAR-CAU-CAV
3	E	302	7L3	CAQ-CAR-CAU-CAW
3	D	303	7L3	CAS-CAR-CAU-CAV
3	F	302	7L3	CAQ-CAR-CAU-CAV
3	G	303	7L3	CAQ-CAR-CAU-CAW
3	F	302	7L3	CAQ-CAR-CAU-CAX
3	E	302	7L3	CAQ-CAR-CAU-CAV
3	E	302	7L3	CAP-CAN-SAL-OAM
3	G	303	7L3	CAQ-CAR-CAU-CAV
3	A	302	7L3	CAQ-CAR-CAU-CAV
3	F	302	7L3	CAS-CAR-CAU-CAV
3	F	302	7L3	CAS-CAR-CAU-CAX
3	C	303	7L3	CAB-NAK-SAL-OAM
3	F	302	7L3	CAB-NAK-SAL-OAM

There are no ring outliers.

9 monomers are involved in 23 short contacts:

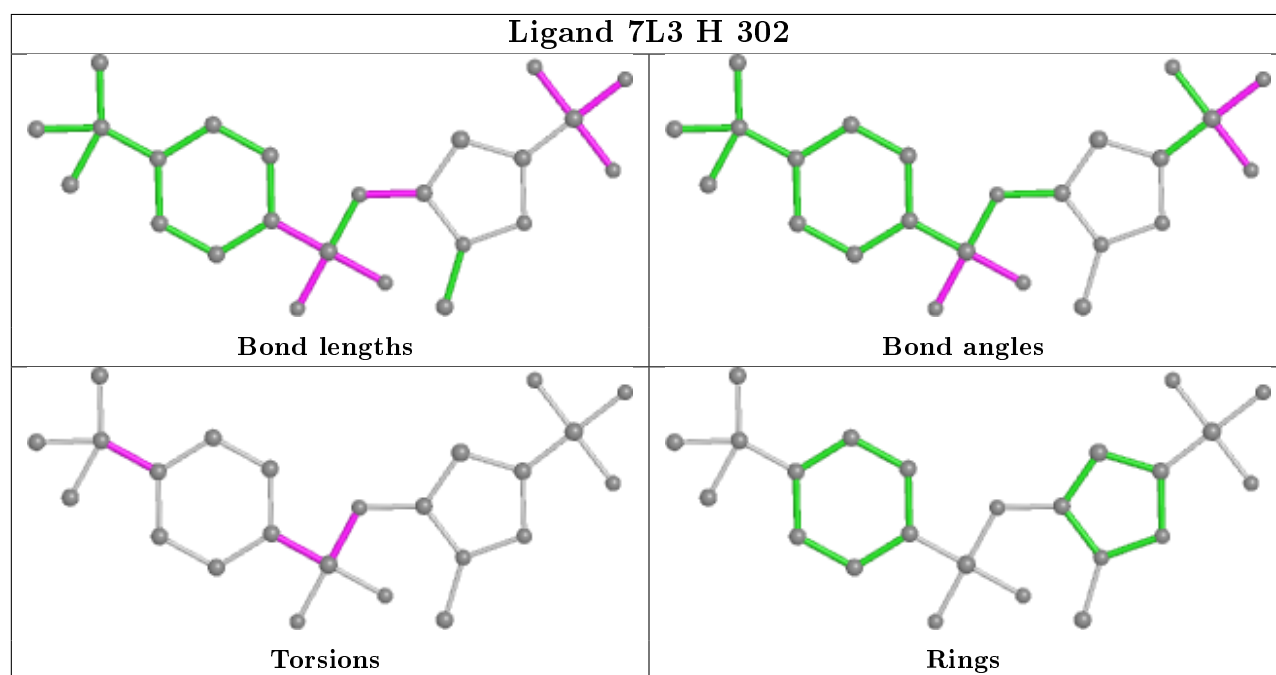
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	302	7L3	2	0
3	D	303	7L3	3	0
3	C	303	7L3	3	0
3	G	303	7L3	2	0
4	C	302	GOL	1	0

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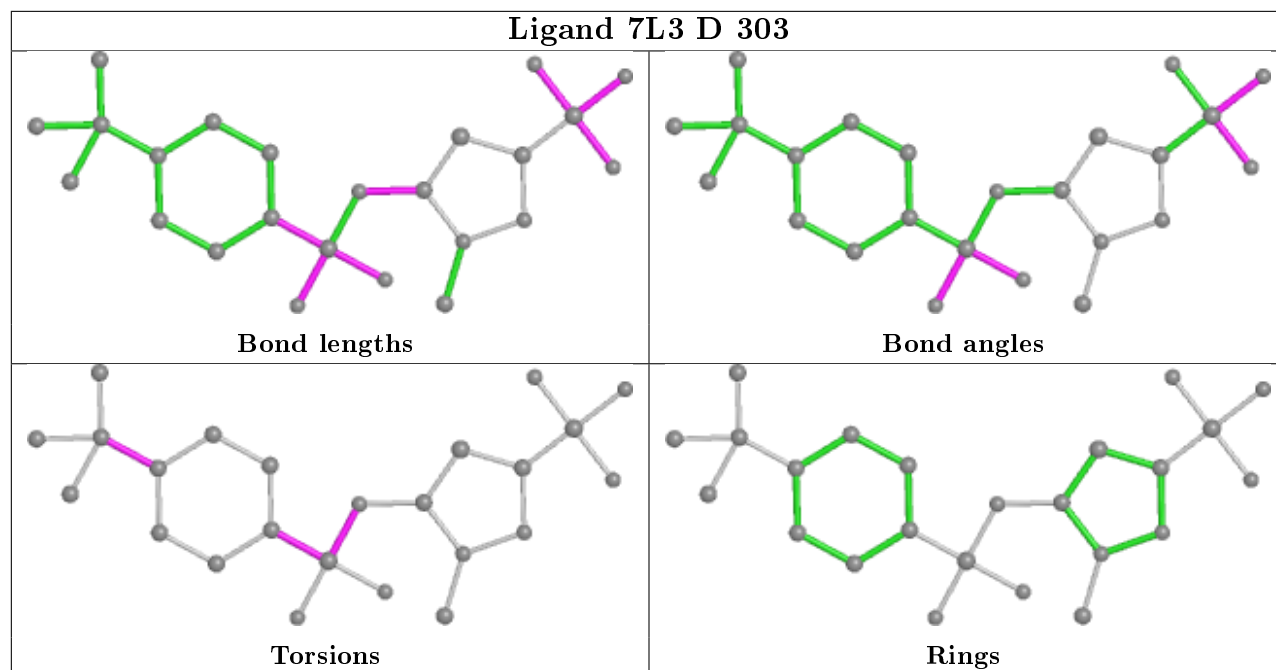
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	7L3	3	0
3	A	302	7L3	3	0
3	F	302	7L3	4	0
3	E	302	7L3	2	0

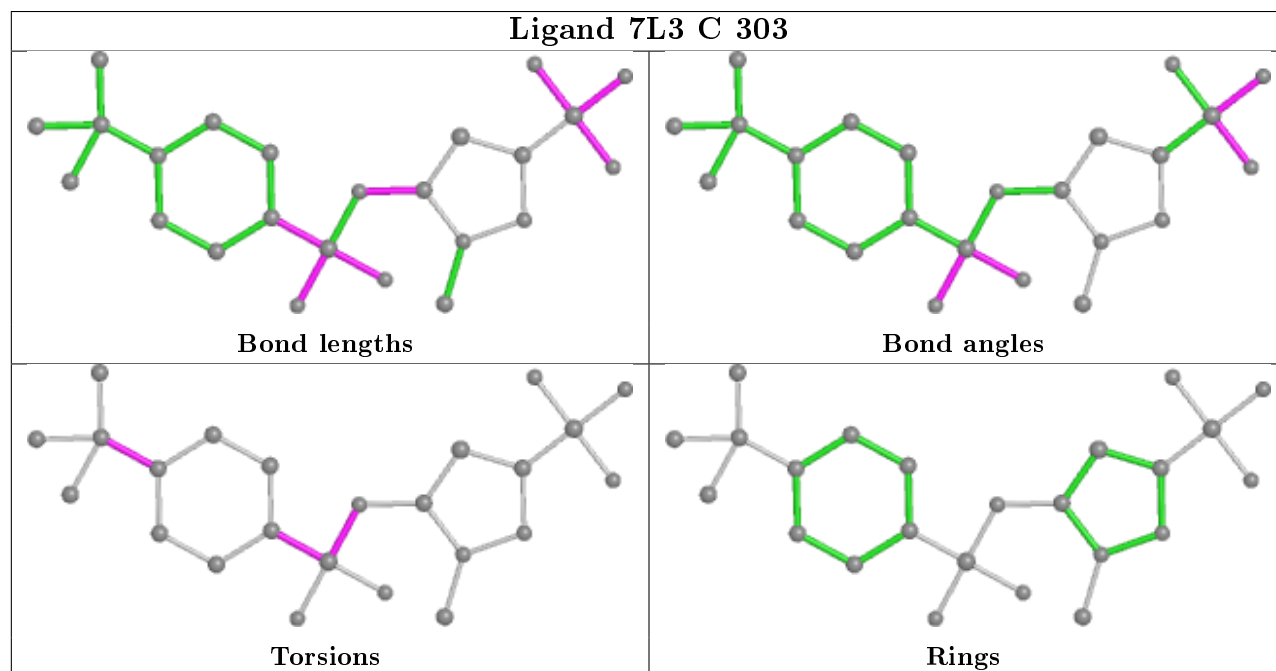
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 7L3 D 303

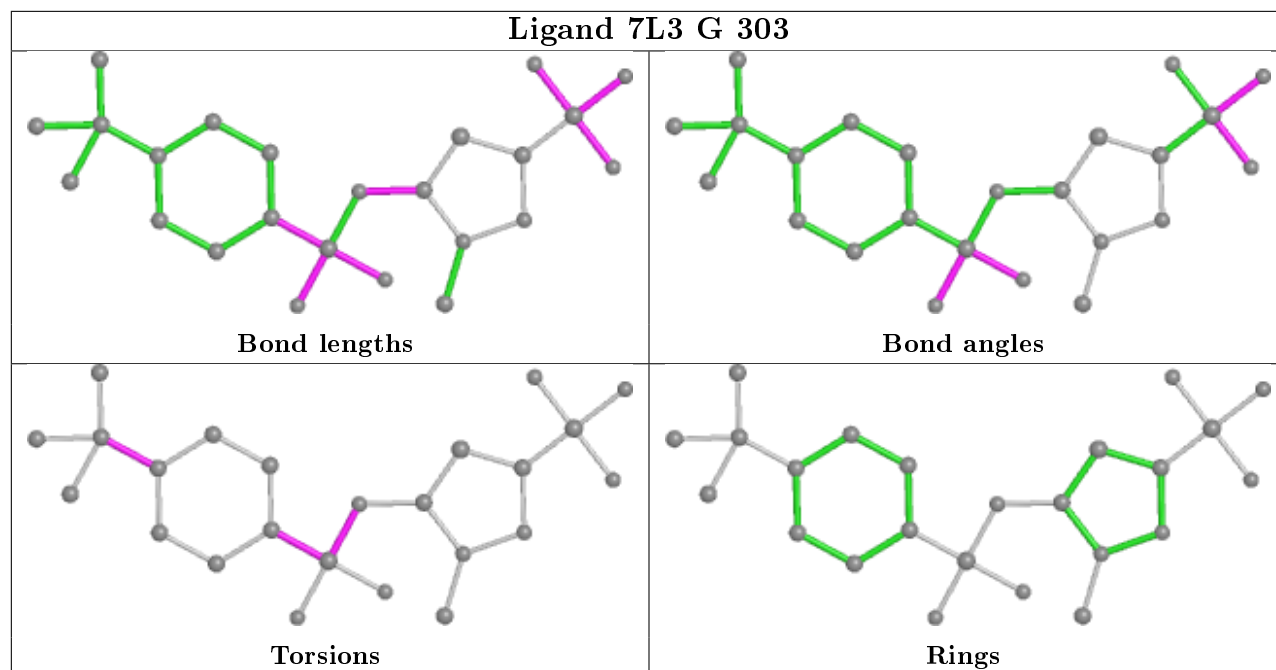


## Ligand 7L3 C 303

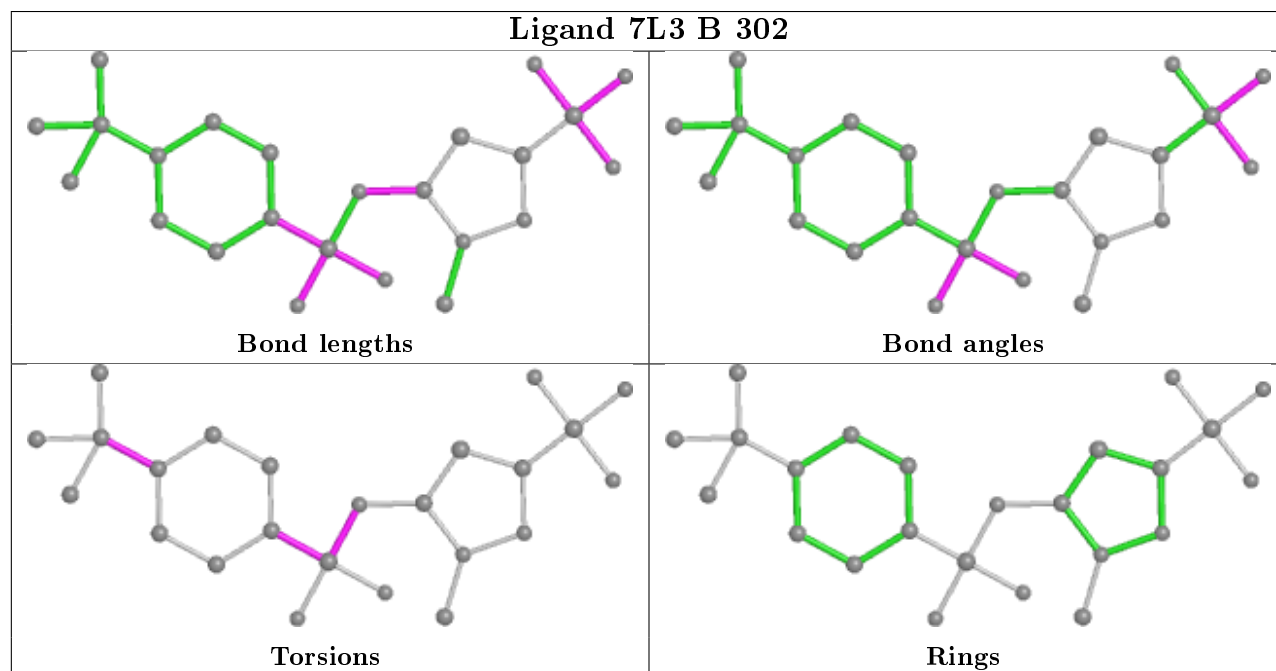




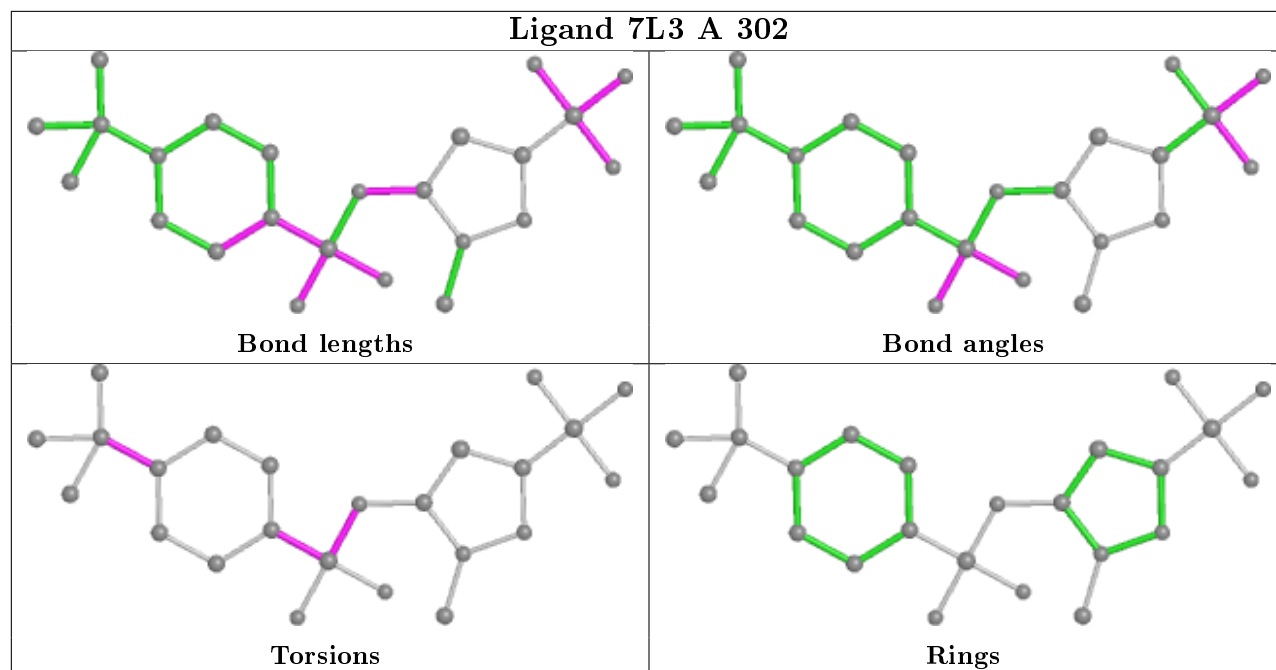
## Ligand 7L3 G 303



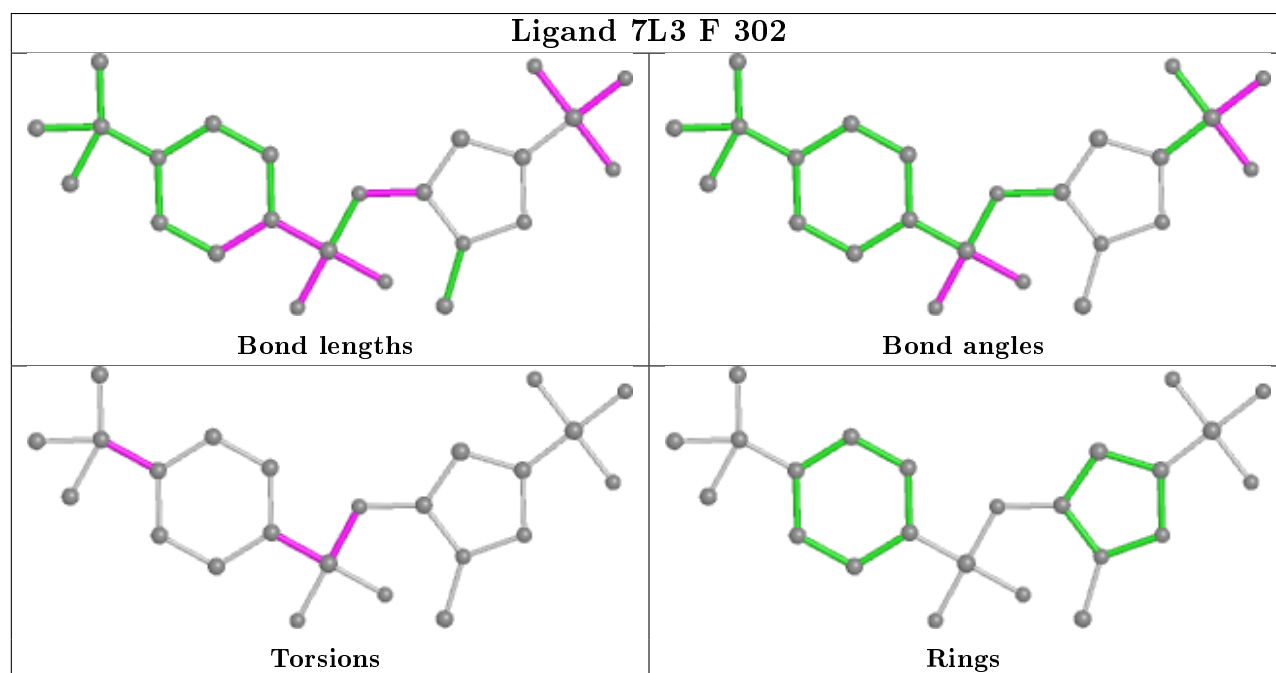
## Ligand 7L3 B 302

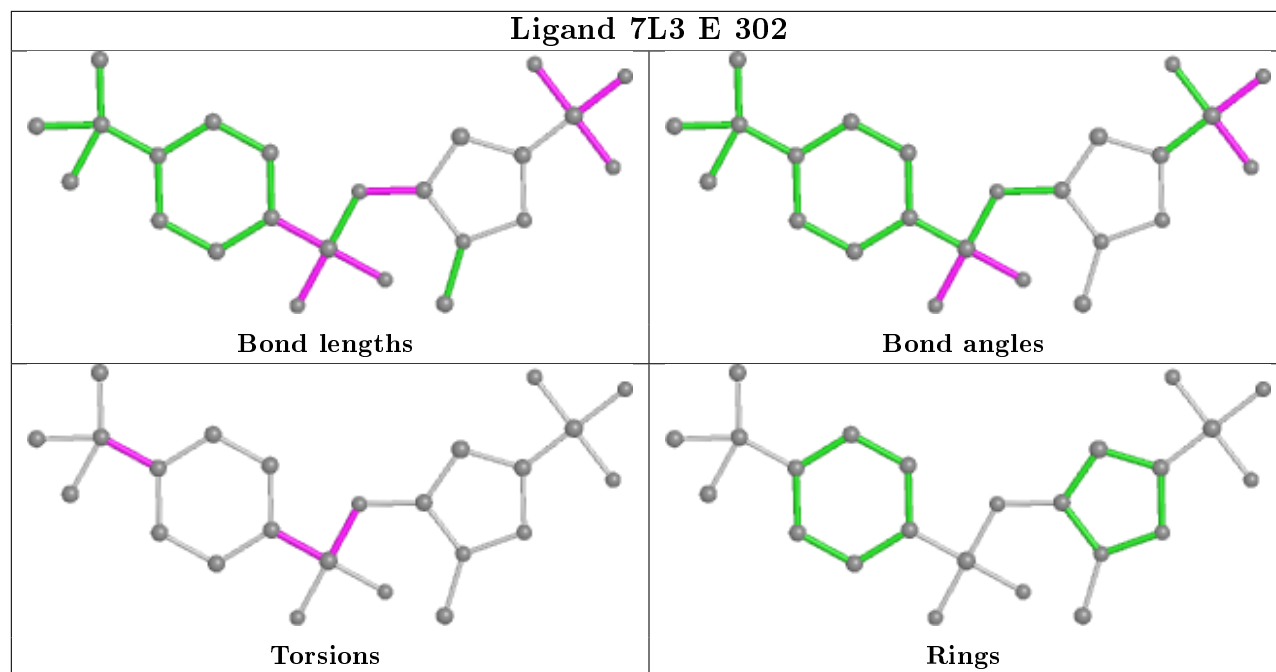


## Ligand 7L3 A 302



## Ligand 7L3 F 302





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	226/234 (96%)	-0.29	1 (0%) 92 93	25, 50, 76, 95	0
1	B	218/234 (93%)	-0.52	0 100 100	21, 39, 60, 83	0
1	C	225/234 (96%)	-0.43	1 (0%) 92 93	18, 37, 56, 88	0
1	D	220/234 (94%)	-0.29	0 100 100	25, 48, 74, 86	0
1	E	226/234 (96%)	-0.47	1 (0%) 92 93	17, 36, 61, 81	0
1	F	222/234 (94%)	-0.35	1 (0%) 91 91	22, 49, 72, 89	0
1	G	222/234 (94%)	-0.31	0 100 100	24, 50, 79, 87	0
1	H	225/234 (96%)	-0.45	0 100 100	20, 40, 62, 78	0
All	All	1784/1872 (95%)	-0.39	4 (0%) 95 95	17, 43, 72, 95	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	135	ALA	2.6
1	A	75	SER	2.4
1	F	27	ASN	2.4
1	E	64	ASP	2.1

### 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 carbohydrates in this entry.

## 6.4 Ligands ⓘ

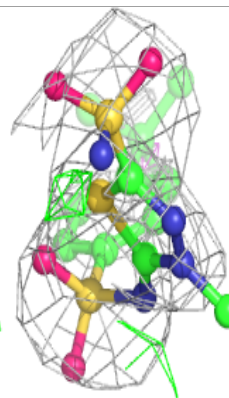
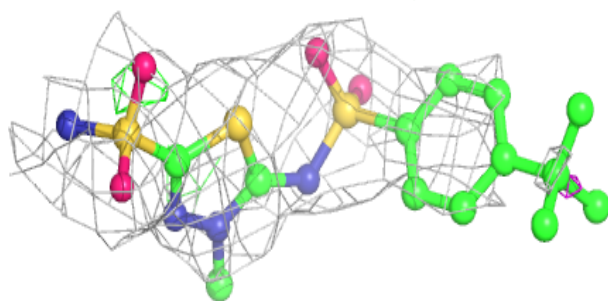
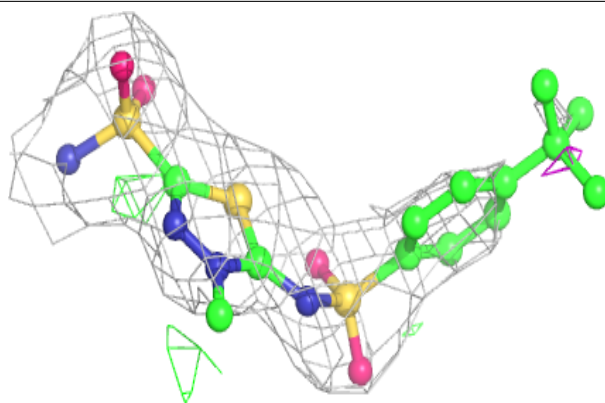
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	CL	G	302	1/1	0.82	0.12	70,70,70,70	0
4	GOL	C	302	6/6	0.91	0.28	42,45,54,64	0
5	CL	D	302	1/1	0.93	0.17	59,59,59,59	0
3	7L3	B	302	24/24	0.94	0.25	32,38,50,59	24
3	7L3	C	303	24/24	0.94	0.23	32,42,50,53	24
3	7L3	F	302	24/24	0.95	0.19	32,40,50,60	24
3	7L3	H	302	24/24	0.95	0.17	32,39,49,52	0
3	7L3	G	303	24/24	0.96	0.18	32,37,47,51	0
3	7L3	A	302	24/24	0.96	0.14	32,35,49,57	0
3	7L3	E	302	24/24	0.97	0.18	21,33,41,48	0
2	ZN	H	301	1/1	0.97	0.09	32,32,32,32	0
3	7L3	D	303	24/24	0.97	0.18	32,42,51,51	0
2	ZN	D	301	1/1	0.98	0.10	50,50,50,50	0
2	ZN	G	301	1/1	0.98	0.10	32,32,32,32	0
2	ZN	F	301	1/1	0.98	0.06	33,33,33,33	1
2	ZN	B	301	1/1	0.98	0.10	33,33,33,33	0
2	ZN	C	301	1/1	0.98	0.09	31,31,31,31	0
2	ZN	A	301	1/1	0.99	0.06	30,30,30,30	1
2	ZN	E	301	1/1	0.99	0.14	32,32,32,32	0

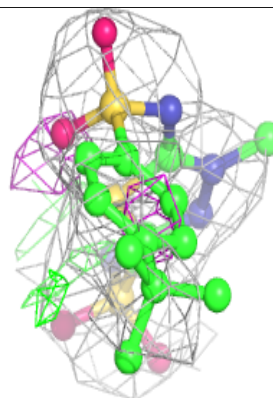
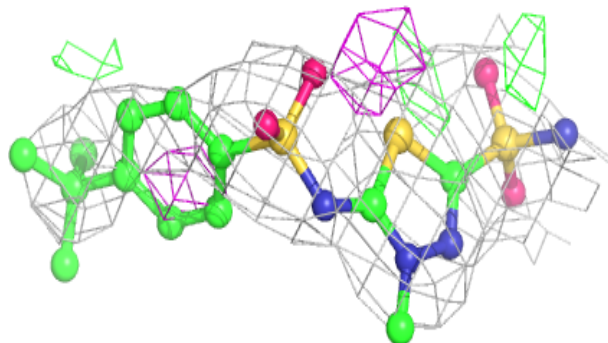
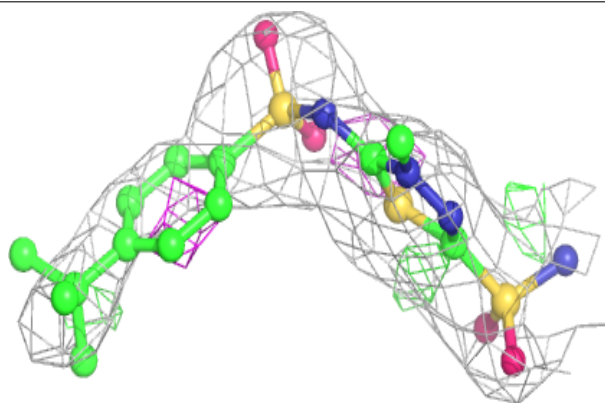
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around 7L3 B 302:**

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

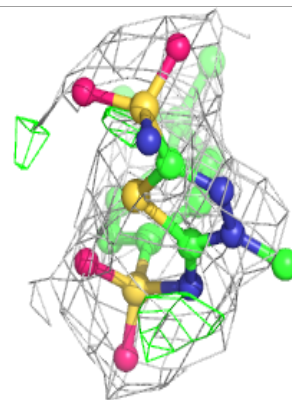
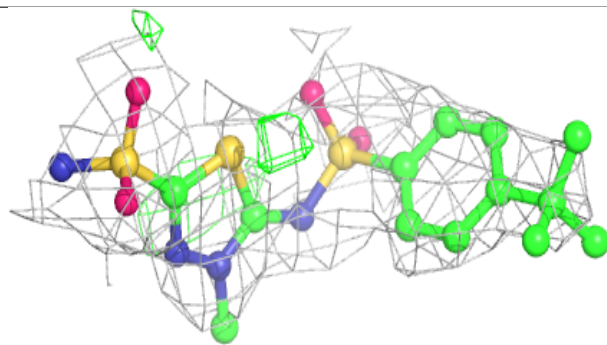
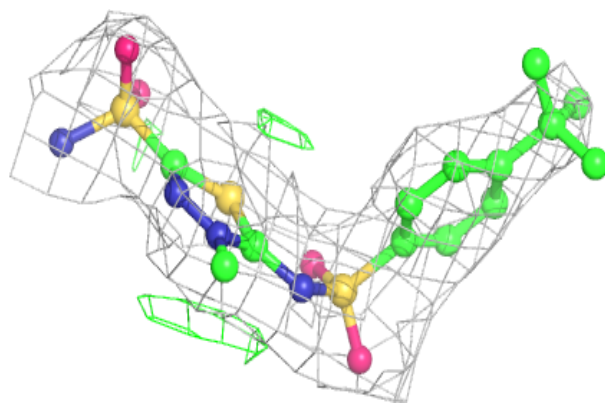
**Electron density around 7L3 C 303:**

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

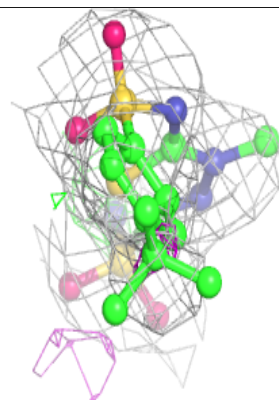
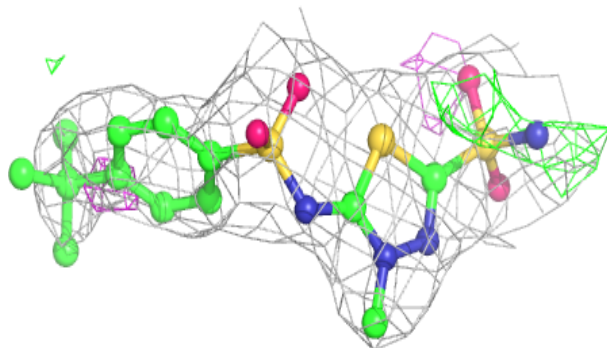
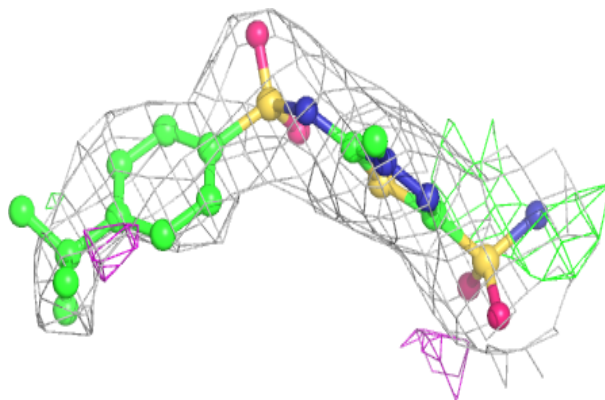


**Electron density around 7L3 F 302:**

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

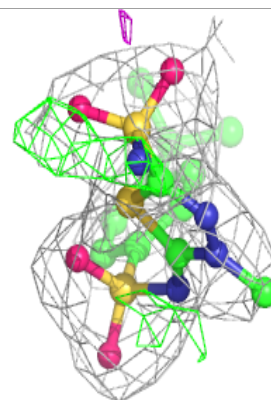
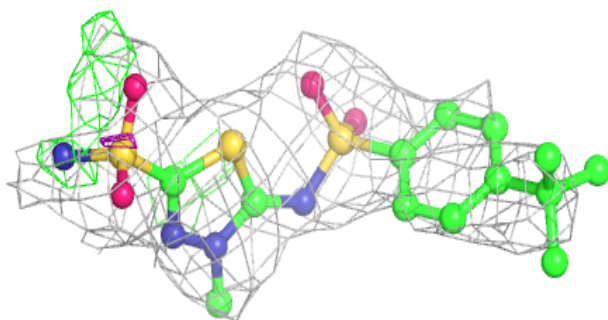
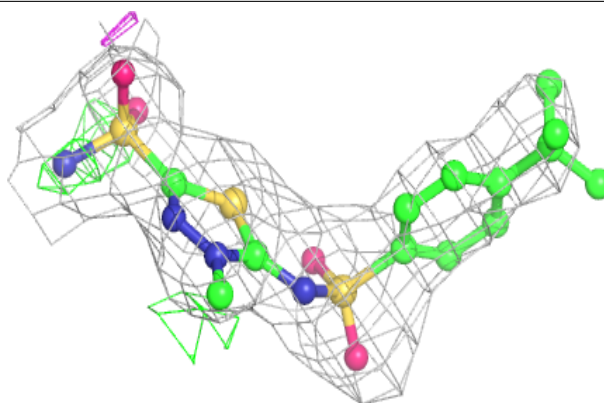
**Electron density around 7L3 H 302:**

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

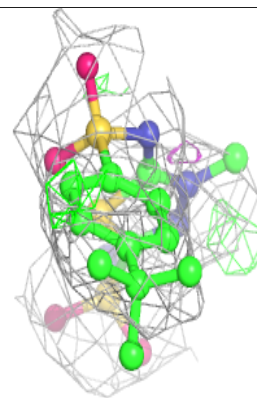
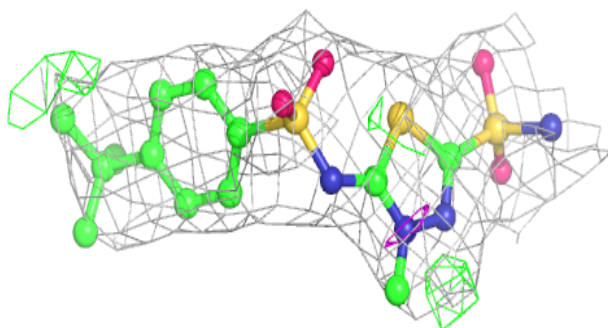
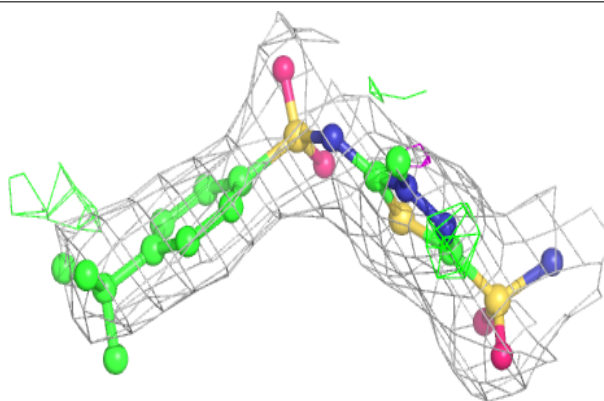


**Electron density around 7L3 G 303:**

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

**Electron density around 7L3 A 302:**

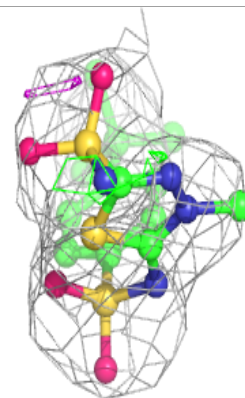
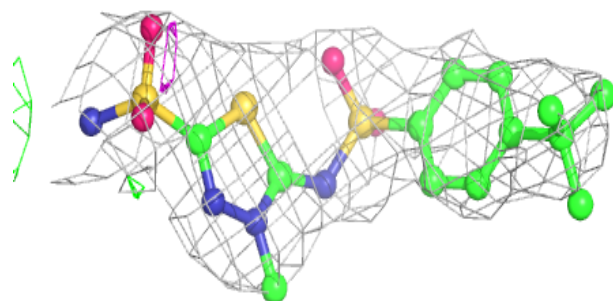
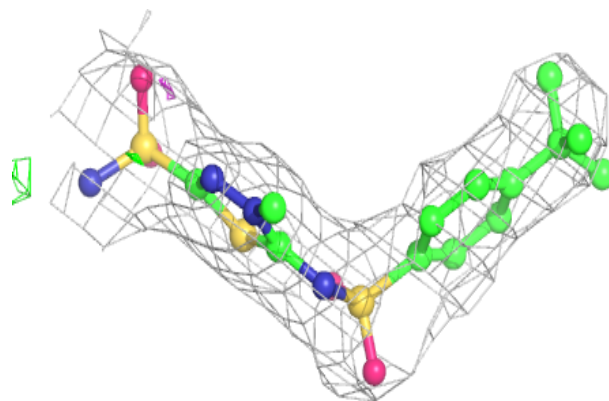
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



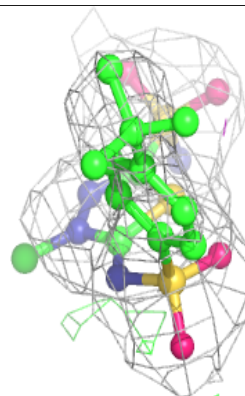
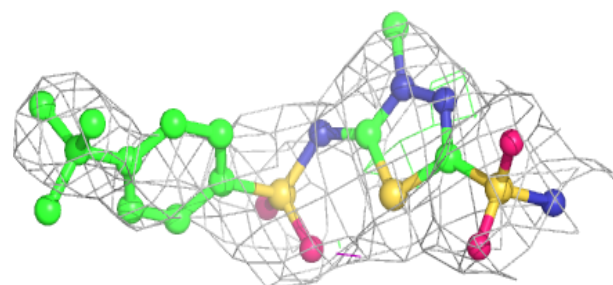
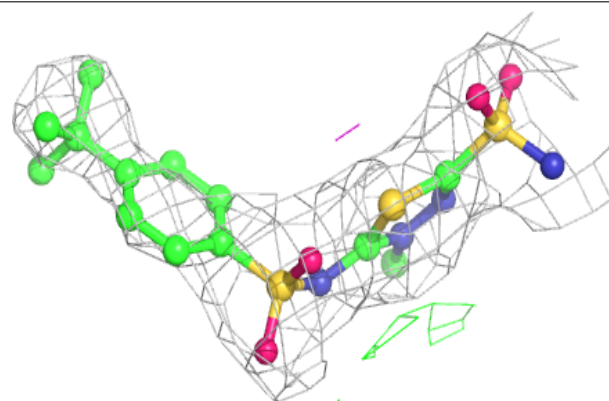


**Electron density around 7L3 E 302:**

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

**Electron density around 7L3 D 303:**

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



## 6.5 Other polymers

There are no such residues in this entry.