



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 10, 2023 – 12:38 am GMT

PDB ID : 8CO0  
Title : Three dimensional structure of human carbonic anhydrase IX in complex with sulfonamide  
Authors : Leitans, J.; Tars, K.  
Deposited on : 2023-02-25  
Resolution : 2.30 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.

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 : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

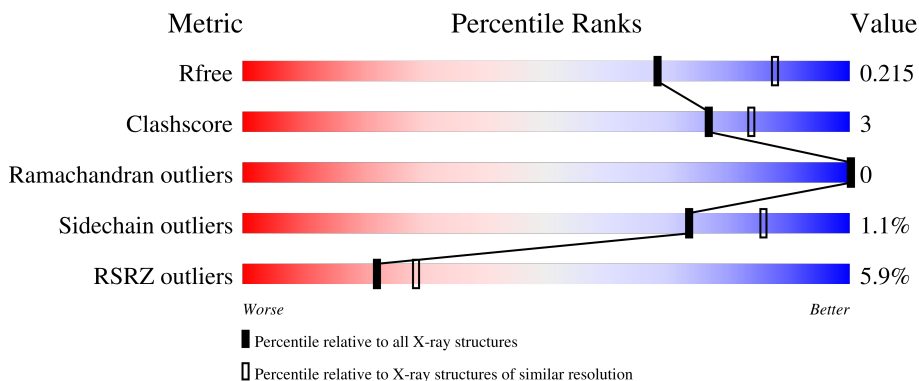
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 2.30 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))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	257	 2% 89% 8% ..
1	C	257	 7% 90% 7% .
1	E	257	 7% 88% 8% ..
1	G	257	 6% 86% 12% .

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
4	GOL	G	303	-	-	X	-

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 8493 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 Carbonic anhydrase 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	251	1946	1240	346	356	4	0	0	0
1	C	251	1940	1237	343	356	4	0	0	0
1	E	251	1946	1240	346	356	4	0	0	0
1	G	251	1946	1240	346	356	4	0	0	0

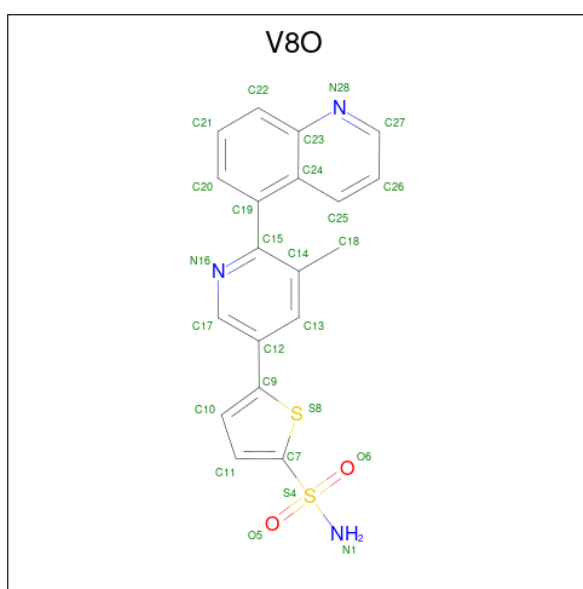
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	GLY	-	expression tag	UNP Q16790
A	4	PRO	-	expression tag	UNP Q16790
A	42	SER	CYS	engineered mutation	UNP Q16790
A	214	GLN	ASN	engineered mutation	UNP Q16790
C	3	GLY	-	expression tag	UNP Q16790
C	4	PRO	-	expression tag	UNP Q16790
C	42	SER	CYS	engineered mutation	UNP Q16790
C	214	GLN	ASN	engineered mutation	UNP Q16790
E	3	GLY	-	expression tag	UNP Q16790
E	4	PRO	-	expression tag	UNP Q16790
E	42	SER	CYS	engineered mutation	UNP Q16790
E	214	GLN	ASN	engineered mutation	UNP Q16790
G	3	GLY	-	expression tag	UNP Q16790
G	4	PRO	-	expression tag	UNP Q16790
G	42	SER	CYS	engineered mutation	UNP Q16790
G	214	GLN	ASN	engineered mutation	UNP Q16790

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0

- Molecule 3 is 5-(5-methyl-6-quinolin-5-yl-pyridin-3-yl)thiophene-2-sulfonamide (three-letter code: V8O) (formula: C<sub>19</sub>H<sub>15</sub>N<sub>3</sub>O<sub>2</sub>S<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O S 26 19 3 2 2	0	0
3	C	1	Total C N O S 26 19 3 2 2	0	0
3	E	1	Total C N O S 26 19 3 2 2	0	0
3	G	1	Total C N O S 26 19 3 2 2	0	0

- 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	A	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	E	1	Total C O 6 3 3	0	0
4	G	1	Total C O 6 3 3	0	0

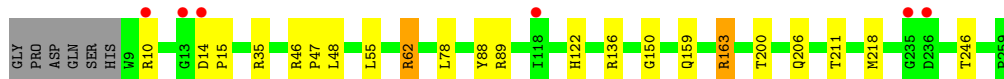
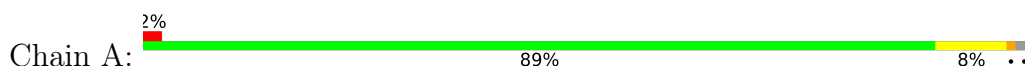
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	192	Total O 192 192	0	0
5	C	136	Total O 136 136	0	0
5	E	132	Total O 132 132	0	0
5	G	123	Total O 123 123	0	0

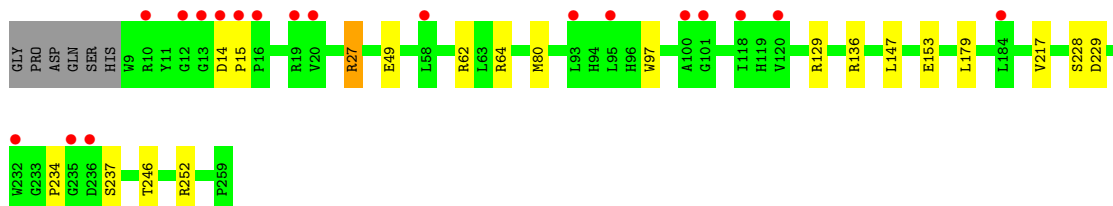
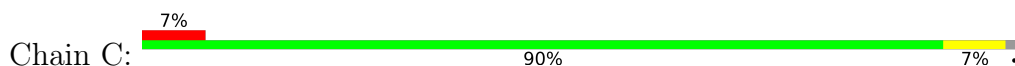
### 3 Residue-property plots [i](#)

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.

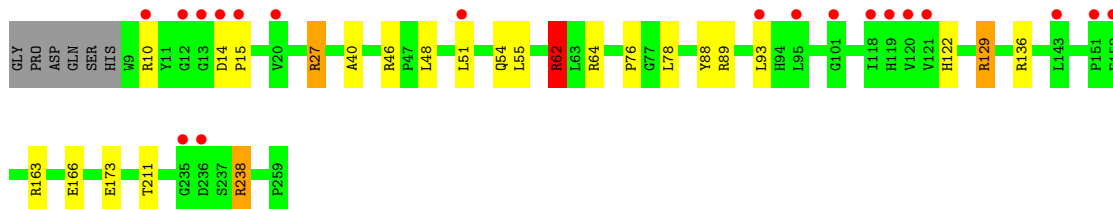
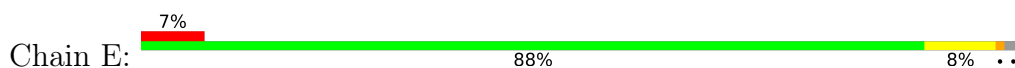
- Molecule 1: Carbonic anhydrase 9



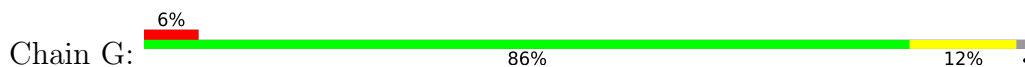
- Molecule 1: Carbonic anhydrase 9

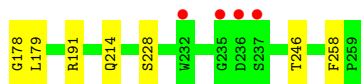


- Molecule 1: Carbonic anhydrase 9



- Molecule 1: Carbonic anhydrase 9







## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	152.95Å 152.95Å 172.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.93 – 2.30 29.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.93-2.30) 100.0 (29.93-2.30)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.31 (at 2.31Å)	Xtrriage
Refinement program	REFMAC 5.8.0350	Depositor
R, $R_{free}$	0.160 , 0.212 0.169 , 0.215	Depositor DCC
$R_{free}$ test set	3427 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.8	Xtrriage
Anisotropy	0.032	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 42.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.007 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8493	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, GOL, V8O

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	0.50	0/2004	0.76	1/2736 (0.0%)
1	C	0.47	0/1998	0.74	0/2729
1	E	0.42	0/2004	0.75	3/2736 (0.1%)
1	G	0.43	0/2004	0.71	0/2736
All	All	0.45	0/8010	0.74	4/10937 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	C	0	6
1	E	0	8
1	G	0	9
All	All	0	28

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	62	ARG	NE-CZ-NH2	-6.69	116.95	120.30
1	E	27	ARG	NE-CZ-NH1	6.62	123.61	120.30
1	A	62	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	E	27	ARG	NE-CZ-NH2	-5.32	117.64	120.30

There are no chirality outliers.

All (28) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	ARG	Sidechain
1	A	136	ARG	Sidechain
1	A	163	ARG	Sidechain
1	A	35	ARG	Sidechain
1	A	62	ARG	Sidechain
1	C	129	ARG	Sidechain
1	C	136	ARG	Sidechain
1	C	252	ARG	Sidechain
1	C	27	ARG	Sidechain
1	C	62	ARG	Sidechain
1	C	64	ARG	Sidechain
1	E	10	ARG	Sidechain
1	E	129	ARG	Sidechain
1	E	136	ARG	Sidechain
1	E	238	ARG	Sidechain
1	E	27	ARG	Sidechain
1	E	46	ARG	Sidechain
1	E	62	ARG	Sidechain
1	E	64	ARG	Sidechain
1	G	10	ARG	Sidechain
1	G	102	ARG	Sidechain
1	G	113	ARG	Sidechain
1	G	129	ARG	Sidechain
1	G	191	ARG	Sidechain
1	G	27	ARG	Sidechain
1	G	35	ARG	Sidechain
1	G	64	ARG	Sidechain
1	G	86	ARG	Sidechain

## 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	1946	0	1899	10	0
1	C	1940	0	1888	8	0
1	E	1946	0	1899	15	0
1	G	1946	0	1899	15	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0

*Continued on next page...*

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1	0	0	0	0
2	G	1	0	0	0	0
3	A	26	0	0	2	0
3	C	26	0	0	0	0
3	E	26	0	0	0	0
3	G	26	0	0	1	0
4	A	6	0	8	0	0
4	C	6	0	8	0	0
4	E	6	0	8	0	0
4	G	6	0	8	4	0
5	A	192	0	0	2	0
5	C	136	0	0	3	0
5	E	132	0	0	1	0
5	G	123	0	0	3	0
All	All	8493	0	7617	47	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 (47) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:GLU:HG2	5:C:522:HOH:O	1.73	0.86
1:G:125:THR:HG21	5:G:448:HOH:O	1.96	0.64
1:G:134:LEU:HD11	3:G:302:V8O:C26	2.29	0.63
1:E:40:ALA:HB1	1:G:43:PRO:HG2	1.83	0.60
1:E:163:ARG:NH2	5:E:401:HOH:O	2.34	0.60
1:C:153:GLU:CG	5:C:522:HOH:O	2.42	0.57
1:C:27:ARG:NH2	5:C:401:HOH:O	2.38	0.56
1:E:93:LEU:C	1:E:93:LEU:HD12	2.26	0.56
4:G:303:GOL:H12	5:G:504:HOH:O	2.06	0.55
1:E:62:ARG:HD3	1:E:173:GLU:OE2	2.08	0.54
1:A:55:LEU:HD21	1:A:78:LEU:HD11	1.88	0.54
1:G:92:GLN:NE2	4:G:303:GOL:O1	2.36	0.54
1:E:14:ASP:HB3	1:E:15:PRO:HA	1.90	0.54
1:C:14:ASP:HB3	1:C:15:PRO:HA	1.93	0.50
1:A:206:GLN:NE2	5:A:406:HOH:O	2.43	0.50
1:E:55:LEU:HD21	1:E:78:LEU:HD11	1.92	0.50
1:A:89:ARG:CZ	1:E:76:PRO:HG2	2.42	0.50
1:G:92:GLN:HE22	4:G:303:GOL:HO1	1.56	0.50
1:A:159:GLN:NE2	5:A:405:HOH:O	2.43	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:THR:OG1	3:A:302:V8O:N1	2.46	0.49
1:E:163:ARG:NH1	1:E:166:GLU:OE2	2.46	0.49
1:E:238:ARG:HH11	1:E:238:ARG:HG2	1.77	0.49
1:G:88:TYR:HB3	1:G:122:HIS:HB3	1.96	0.48
1:A:14:ASP:HB3	1:A:15:PRO:HA	1.96	0.47
1:G:56:PRO:O	1:G:178:GLY:HA3	2.14	0.47
1:A:150:GLY:CA	1:A:218:MET:HE3	2.45	0.46
1:A:88:TYR:HB3	1:A:122:HIS:HB3	1.97	0.46
1:C:49:GLU:O	1:C:80:MET:HA	2.16	0.45
1:E:163:ARG:HD2	1:E:166:GLU:OE1	2.17	0.45
1:E:48:LEU:HD21	1:E:211:THR:HG21	1.98	0.45
1:E:88:TYR:HB3	1:E:122:HIS:HB3	1.99	0.45
1:G:163:ARG:NH2	5:G:410:HOH:O	2.49	0.44
1:G:92:GLN:NE2	4:G:303:GOL:HO1	2.14	0.44
3:A:302:V8O:C27	1:E:129:ARG:HD2	2.47	0.43
1:E:89:ARG:HH11	1:E:89:ARG:HG2	1.83	0.43
1:C:97:TRP:CD1	1:C:228:SER:HA	2.53	0.43
1:C:234:PRO:O	1:C:237:SER:HB3	2.18	0.43
1:G:95:LEU:HD23	1:G:118:ILE:HA	2.01	0.42
1:C:147:LEU:HD22	1:C:217:VAL:CG2	2.49	0.42
1:A:48:LEU:HD21	1:A:211:THR:HG21	2.02	0.42
1:G:97:TRP:CD1	1:G:228:SER:HA	2.55	0.41
1:G:49:GLU:C	1:G:50:LEU:HD12	2.40	0.41
1:E:163:ARG:HH11	1:E:166:GLU:CD	2.23	0.41
1:G:14:ASP:HB3	1:G:15:PRO:HA	2.01	0.41
1:A:46:ARG:HB3	1:A:47:PRO:HD2	2.01	0.41
1:G:37:GLN:HE21	1:G:37:GLN:HB2	1.58	0.41
1:G:214:GLN:HG3	1:G:258:PHE:CZ	2.56	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	249/257 (97%)	245 (98%)	4 (2%)	0	100	100
1	C	249/257 (97%)	245 (98%)	4 (2%)	0	100	100
1	E	249/257 (97%)	242 (97%)	7 (3%)	0	100	100
1	G	249/257 (97%)	243 (98%)	6 (2%)	0	100	100
All	All	996/1028 (97%)	975 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	204/210 (97%)	202 (99%)	2 (1%)	76	87
1	C	203/210 (97%)	200 (98%)	3 (2%)	65	79
1	E	204/210 (97%)	202 (99%)	2 (1%)	76	87
1	G	204/210 (97%)	202 (99%)	2 (1%)	76	87
All	All	815/840 (97%)	806 (99%)	9 (1%)	73	86

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	ARG
1	A	246	THR
1	C	179	LEU
1	C	229	ASP
1	C	246	THR
1	E	51	LEU
1	E	54	GLN
1	G	179	LEU
1	G	246	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	159	GLN
1	G	37	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
4	GOL	A	303	-	5,5,5	0.13	0	5,5,5	0.38	0
3	V8O	A	302	2	25,29,29	2.90	11 (44%)	32,43,43	5.15	20 (62%)
4	GOL	C	303	-	5,5,5	0.16	0	5,5,5	0.37	0
4	GOL	E	303	-	5,5,5	0.13	0	5,5,5	0.27	0
3	V8O	E	302	2	25,29,29	2.59	11 (44%)	32,43,43	4.27	16 (50%)
3	V8O	C	302	2	25,29,29	2.57	5 (20%)	32,43,43	5.27	22 (68%)
4	GOL	G	303	-	5,5,5	0.08	0	5,5,5	0.35	0
3	V8O	G	302	2	25,29,29	2.09	5 (20%)	32,43,43	4.42	15 (46%)

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	GOL	A	303	-	-	2/4/4/4	-
3	V8O	A	302	2	-	0/8/14/14	0/4/4/4
4	GOL	C	303	-	-	2/4/4/4	-
4	GOL	E	303	-	-	4/4/4/4	-
3	V8O	E	302	2	-	0/8/14/14	0/4/4/4
3	V8O	C	302	2	-	0/8/14/14	0/4/4/4
4	GOL	G	303	-	-	4/4/4/4	-
3	V8O	G	302	2	-	0/8/14/14	0/4/4/4

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	302	V8O	C15-C14	9.15	1.52	1.40
3	A	302	V8O	C15-C14	8.44	1.51	1.40
3	E	302	V8O	C15-C14	7.96	1.50	1.40
3	G	302	V8O	C15-C14	5.67	1.47	1.40
3	C	302	V8O	C24-C23	5.48	1.51	1.42
3	G	302	V8O	C7-S4	-5.22	1.68	1.76
3	A	302	V8O	C24-C23	4.95	1.50	1.42
3	A	302	V8O	C7-S8	4.71	1.80	1.72
3	G	302	V8O	C24-C23	4.47	1.49	1.42
3	E	302	V8O	C24-C23	4.44	1.49	1.42
3	C	302	V8O	C7-S4	-4.38	1.69	1.76
3	E	302	V8O	C7-S4	-4.33	1.69	1.76
3	E	302	V8O	C19-C24	3.89	1.51	1.42
3	A	302	V8O	C7-S4	-3.57	1.71	1.76
3	A	302	V8O	C20-C19	3.43	1.44	1.38
3	A	302	V8O	C13-C14	-3.27	1.34	1.39
3	A	302	V8O	C10-C11	-3.10	1.23	1.39
3	C	302	V8O	C20-C19	3.00	1.43	1.38
3	E	302	V8O	C7-S8	2.95	1.77	1.72
3	A	302	V8O	S4-N1	2.94	1.66	1.60
3	E	302	V8O	C26-C25	2.86	1.43	1.36
3	A	302	V8O	O6-S4	2.80	1.49	1.43
3	A	302	V8O	C27-N28	2.78	1.38	1.32
3	A	302	V8O	C19-C24	2.58	1.48	1.42
3	G	302	V8O	C20-C19	2.54	1.42	1.38
3	E	302	V8O	C20-C19	2.46	1.42	1.38
3	E	302	V8O	C27-N28	2.30	1.37	1.32
3	E	302	V8O	O6-S4	2.21	1.47	1.43
3	G	302	V8O	C21-C22	2.17	1.41	1.36

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	302	V8O	C21-C22	2.17	1.41	1.36
3	E	302	V8O	O5-S4	2.07	1.47	1.43
3	C	302	V8O	C27-N28	2.01	1.36	1.32

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	302	V8O	O5-S4-N1	17.56	133.39	107.36
3	G	302	V8O	O5-S4-O6	-14.59	94.78	118.76
3	A	302	V8O	O5-S4-O6	-13.41	96.72	118.76
3	E	302	V8O	O6-S4-N1	13.18	126.91	107.36
3	C	302	V8O	O6-S4-C7	-13.12	88.05	107.29
3	A	302	V8O	O6-S4-N1	12.97	126.59	107.36
3	G	302	V8O	O6-S4-N1	12.83	126.39	107.36
3	C	302	V8O	O5-S4-O6	-11.93	99.15	118.76
3	E	302	V8O	O5-S4-O6	-11.23	100.31	118.76
3	A	302	V8O	O5-S4-C7	-11.10	91.02	107.29
3	E	302	V8O	O5-S4-C7	-10.43	91.99	107.29
3	A	302	V8O	C24-C23-N28	-8.64	113.63	122.83
3	A	302	V8O	C14-C15-N16	-7.38	116.65	123.05
3	C	302	V8O	C24-C23-N28	-7.31	115.04	122.83
3	G	302	V8O	O5-S4-C7	-6.33	98.00	107.29
3	G	302	V8O	C14-C15-N16	-6.29	117.59	123.05
3	A	302	V8O	C13-C12-C9	-6.23	109.97	120.21
3	A	302	V8O	C17-C12-C9	5.86	130.66	121.22
3	C	302	V8O	C14-C15-N16	-5.86	117.97	123.05
3	G	302	V8O	C24-C23-N28	-5.62	116.85	122.83
3	A	302	V8O	C7-S4-N1	5.60	118.60	108.28
3	G	302	V8O	C7-S4-N1	5.16	117.79	108.28
3	E	302	V8O	C24-C23-N28	-5.12	117.38	122.83
3	E	302	V8O	C27-N28-C23	5.06	124.77	116.93
3	A	302	V8O	C18-C14-C15	5.02	129.72	122.77
3	E	302	V8O	O5-S4-N1	4.93	114.67	107.36
3	A	302	V8O	C27-N28-C23	4.77	124.32	116.93
3	E	302	V8O	C14-C15-N16	-4.53	119.12	123.05
3	C	302	V8O	C25-C24-C23	4.52	123.38	118.33
3	C	302	V8O	C22-C23-C24	4.49	124.28	119.13
3	A	302	V8O	C18-C14-C13	-4.34	111.47	119.49
3	C	302	V8O	C18-C14-C15	4.08	128.42	122.77
3	C	302	V8O	C19-C15-C14	4.03	127.24	121.98
3	A	302	V8O	C22-C23-N28	3.92	124.95	118.52
3	C	302	V8O	C7-S4-N1	3.74	115.18	108.28

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	302	V80	O5-S4-N1	-3.65	101.95	107.36
3	G	302	V80	C19-C15-N16	3.65	122.69	115.52
3	C	302	V80	C18-C14-C13	-3.27	113.44	119.49
3	E	302	V80	C22-C23-N28	3.23	123.82	118.52
3	A	302	V80	C19-C24-C23	-3.16	115.73	117.44
3	C	302	V80	C27-N28-C23	3.16	121.82	116.93
3	G	302	V80	C26-C25-C24	-3.16	116.51	120.89
3	G	302	V80	C25-C24-C23	3.11	121.81	118.33
3	E	302	V80	C18-C14-C15	3.11	127.08	122.77
3	G	302	V80	C13-C12-C17	3.11	120.19	117.11
3	E	302	V80	C26-C27-N28	-3.09	119.20	123.94
3	C	302	V80	C26-C25-C24	-3.03	116.69	120.89
3	G	302	V80	C27-N28-C23	2.95	121.50	116.93
3	A	302	V80	C17-N16-C15	2.94	123.27	117.95
3	G	302	V80	C17-N16-C15	2.86	123.12	117.95
3	A	302	V80	C19-C15-N16	2.85	121.13	115.52
3	C	302	V80	C13-C12-C17	2.77	119.85	117.11
3	C	302	V80	C21-C20-C19	2.70	126.24	121.09
3	A	302	V80	C13-C14-C15	2.62	118.79	117.12
3	C	302	V80	C17-C12-C9	-2.60	117.03	121.22
3	E	302	V80	C19-C24-C23	-2.56	116.05	117.44
3	E	302	V80	C13-C12-C17	2.54	119.63	117.11
3	A	302	V80	C25-C24-C23	2.54	121.17	118.33
3	G	302	V80	C12-C17-N16	-2.43	120.32	124.32
3	E	302	V80	C18-C14-C13	-2.42	115.02	119.49
3	C	302	V80	C17-N16-C15	2.38	122.25	117.95
3	A	302	V80	C12-C17-N16	-2.37	120.41	124.32
3	E	302	V80	C21-C22-C23	2.33	123.43	120.08
3	C	302	V80	C21-C22-C23	-2.32	116.75	120.08
3	A	302	V80	C13-C12-C17	2.26	119.35	117.11
3	C	302	V80	C25-C24-C19	-2.19	119.23	123.16
3	E	302	V80	C17-N16-C15	2.17	121.88	117.95
3	E	302	V80	C12-C13-C14	-2.17	119.77	122.33
3	C	302	V80	C12-C13-C14	-2.16	119.78	122.33
3	G	302	V80	C22-C23-N28	2.13	122.01	118.52
3	C	302	V80	O5-S4-C7	-2.03	104.30	107.29
3	C	302	V80	C20-C19-C24	-2.03	115.25	118.66
3	A	302	V80	C22-C23-C24	2.00	121.43	119.13

There are no chirality outliers.

All (12) torsion outliers are listed below:

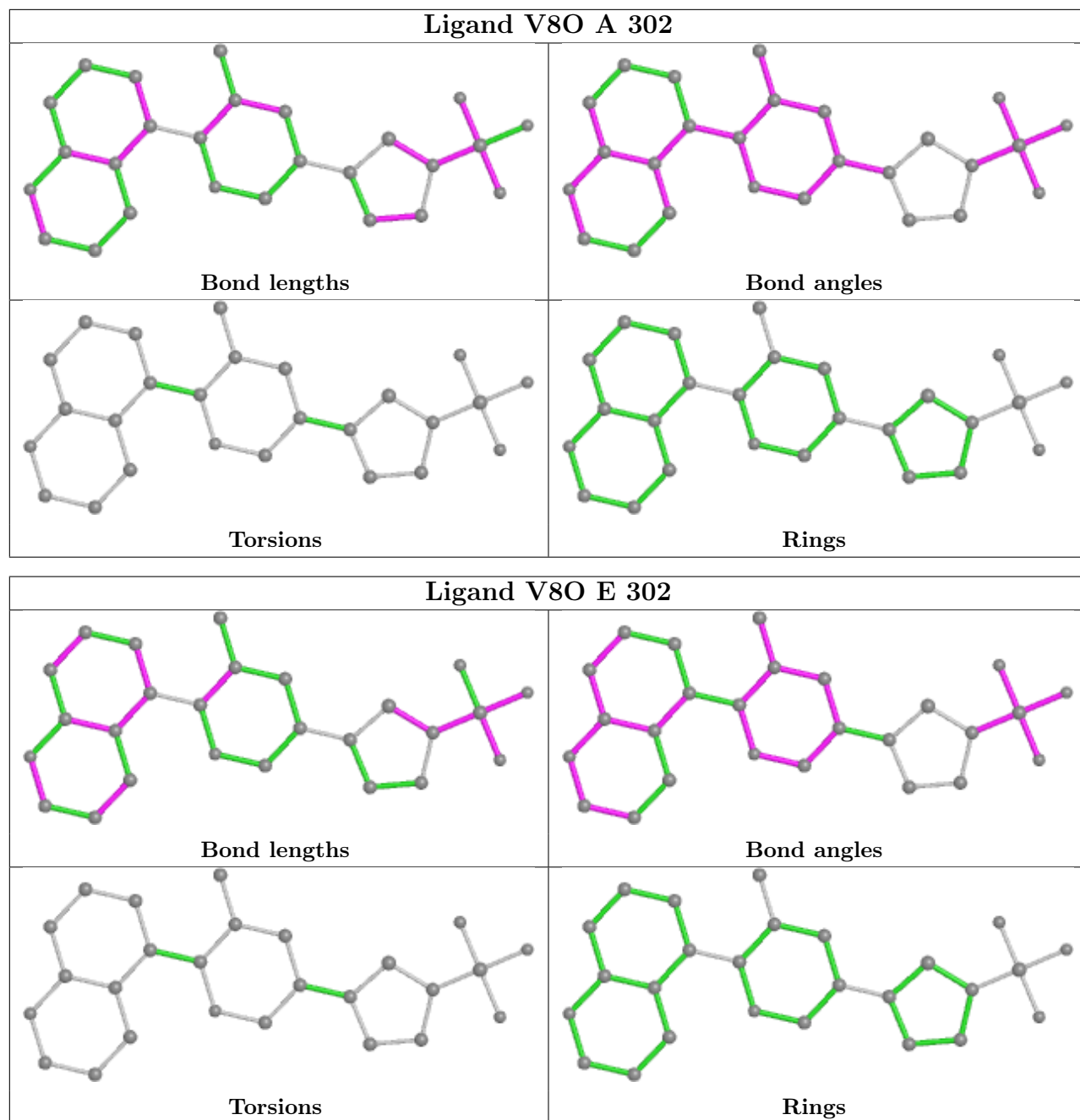
Mol	Chain	Res	Type	Atoms
4	C	303	GOL	O1-C1-C2-C3
4	E	303	GOL	C1-C2-C3-O3
4	G	303	GOL	O1-C1-C2-O2
4	G	303	GOL	O1-C1-C2-C3
4	G	303	GOL	C1-C2-C3-O3
4	A	303	GOL	O1-C1-C2-C3
4	E	303	GOL	O1-C1-C2-C3
4	A	303	GOL	O1-C1-C2-O2
4	C	303	GOL	O1-C1-C2-O2
4	E	303	GOL	O2-C2-C3-O3
4	G	303	GOL	O2-C2-C3-O3
4	E	303	GOL	O1-C1-C2-O2

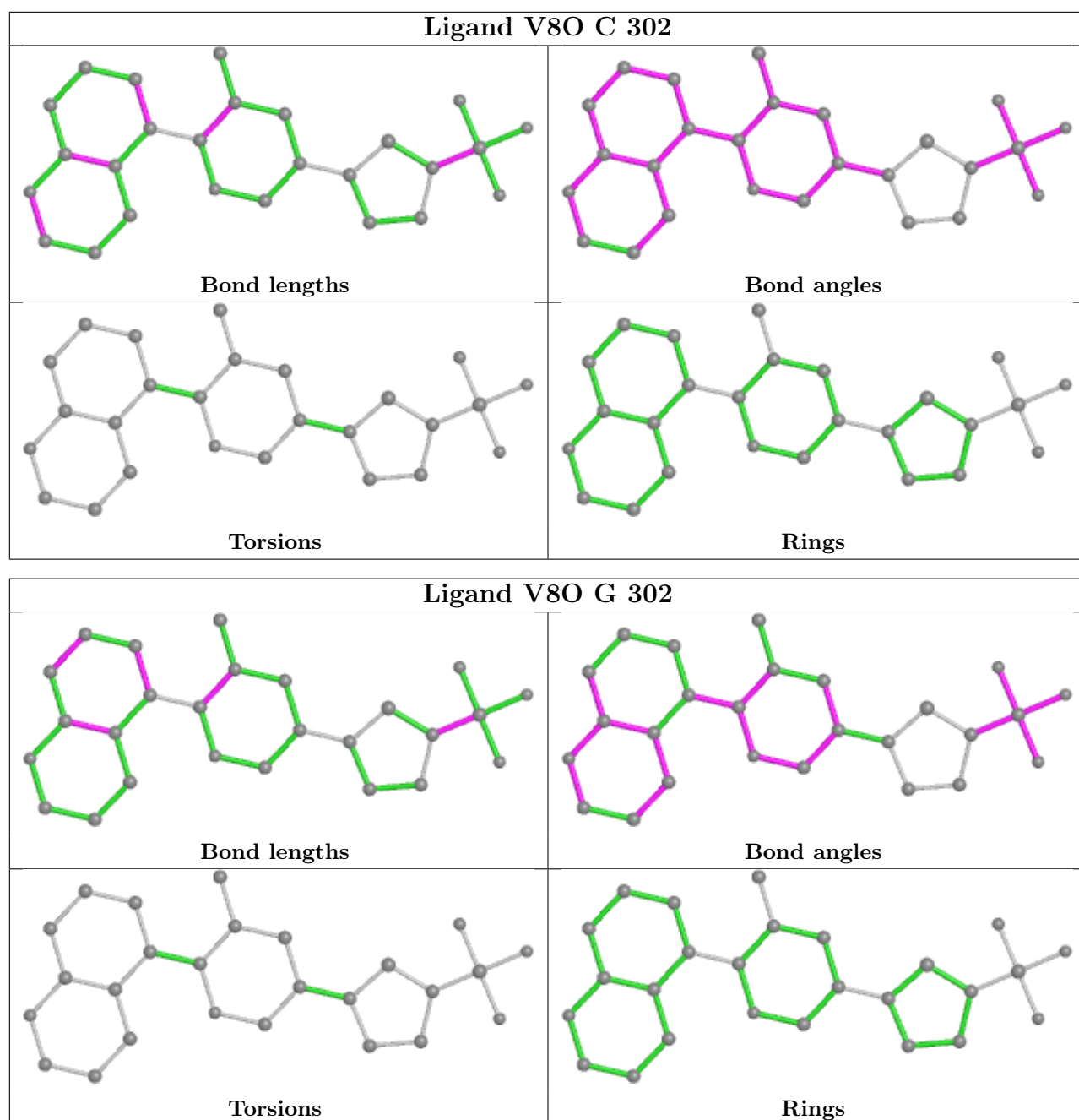
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	V8O	2	0
4	G	303	GOL	4	0
3	G	302	V8O	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, 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	251/257 (97%)	-0.44	6 (2%) 59 66	25, 36, 61, 85	0
1	C	251/257 (97%)	-0.12	19 (7%) 13 18	27, 39, 79, 121	0
1	E	251/257 (97%)	-0.08	19 (7%) 13 18	30, 46, 75, 91	0
1	G	251/257 (97%)	-0.01	15 (5%) 21 28	32, 48, 82, 98	0
All	All	1004/1028 (97%)	-0.16	59 (5%) 22 28	25, 42, 75, 121	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	16	PRO	6.1
1	C	12	GLY	5.9
1	C	13	GLY	5.4
1	G	235	GLY	5.2
1	C	14	ASP	4.7
1	G	13	GLY	4.1
1	G	232	TRP	3.9
1	G	14	ASP	3.8
1	G	236	ASP	3.7
1	E	12	GLY	3.7
1	E	13	GLY	3.7
1	E	118	ILE	3.6
1	G	19	ARG	3.5
1	E	95	LEU	3.4
1	C	93	LEU	3.3
1	E	236	ASP	3.3
1	C	235	GLY	3.2
1	A	236	ASP	3.2
1	C	101	GLY	3.2
1	C	15	PRO	3.2
1	C	236	ASP	3.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	E	120	VAL	3.1
1	G	95	LEU	3.1
1	G	118	ILE	3.0
1	C	118	ILE	3.0
1	G	12	GLY	3.0
1	A	235	GLY	3.0
1	E	101	GLY	3.0
1	A	13	GLY	2.9
1	E	15	PRO	2.9
1	G	10	ARG	2.9
1	G	37	GLN	2.8
1	A	118	ILE	2.7
1	E	235	GLY	2.7
1	E	20	VAL	2.7
1	E	119	HIS	2.6
1	E	10	ARG	2.6
1	C	95	LEU	2.6
1	E	151	PRO	2.6
1	E	93	LEU	2.5
1	C	120	VAL	2.5
1	G	15	PRO	2.5
1	E	14	ASP	2.5
1	A	14	ASP	2.4
1	C	10	ARG	2.4
1	C	20	VAL	2.4
1	A	10	ARG	2.3
1	C	100	ALA	2.3
1	C	232	TRP	2.3
1	C	19	ARG	2.3
1	G	237	SER	2.3
1	E	152	GLU	2.2
1	C	184	LEU	2.2
1	C	58	LEU	2.1
1	G	101	GLY	2.1
1	E	51	LEU	2.1
1	E	143	LEU	2.1
1	E	121	VAL	2.0
1	G	151	PRO	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 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.

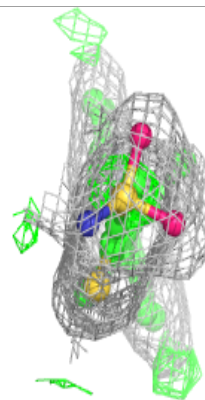
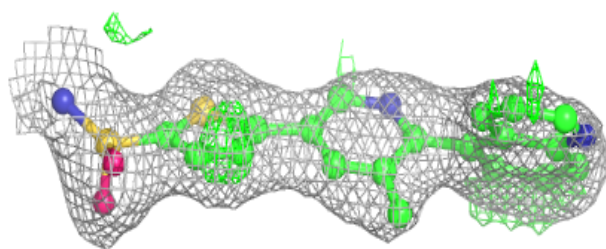
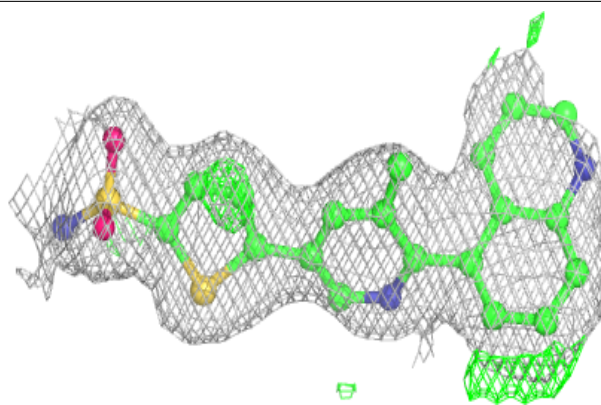
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	GOL	C	303	6/6	0.89	0.28	61,67,70,75	0
4	GOL	E	303	6/6	0.89	0.29	63,74,76,83	0
3	V8O	A	302	26/26	0.94	0.14	25,55,71,76	0
4	GOL	G	303	6/6	0.94	0.32	53,64,66,73	0
4	GOL	A	303	6/6	0.95	0.27	50,58,68,81	0
3	V8O	E	302	26/26	0.95	0.13	32,49,62,65	0
3	V8O	C	302	26/26	0.96	0.13	29,50,61,62	0
3	V8O	G	302	26/26	0.96	0.12	36,51,57,59	0
2	ZN	E	301	1/1	0.99	0.12	36,36,36,36	0
2	ZN	G	301	1/1	0.99	0.11	40,40,40,40	0
2	ZN	A	301	1/1	1.00	0.12	30,30,30,30	0
2	ZN	C	301	1/1	1.00	0.11	33,33,33,33	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.

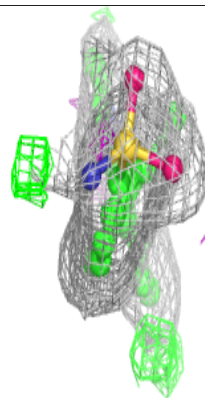
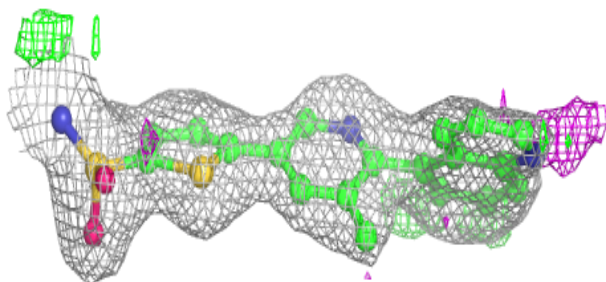
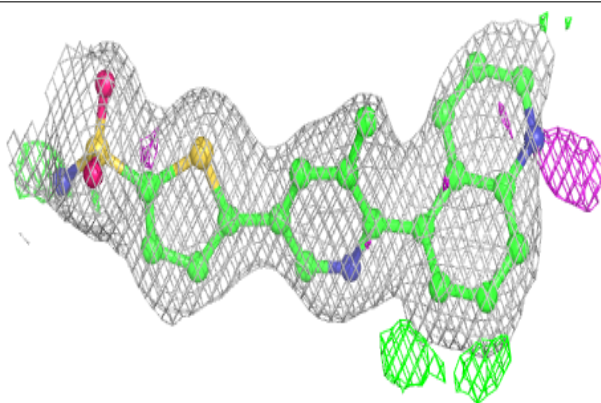


**Electron density around V8O A 302:**

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

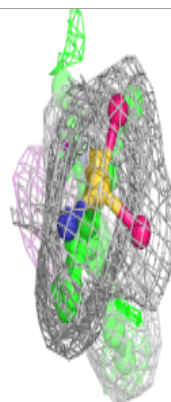
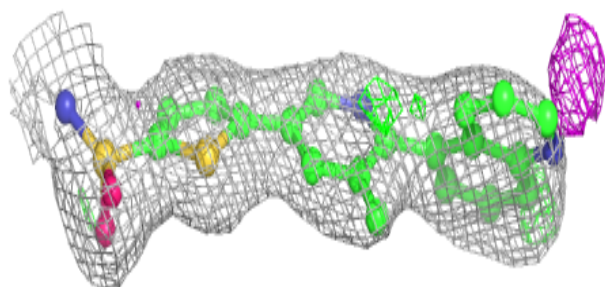
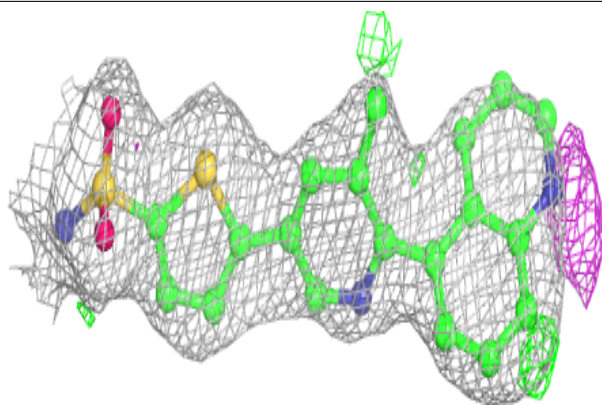
**Electron density around V8O E 302:**

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

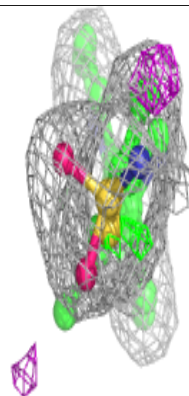
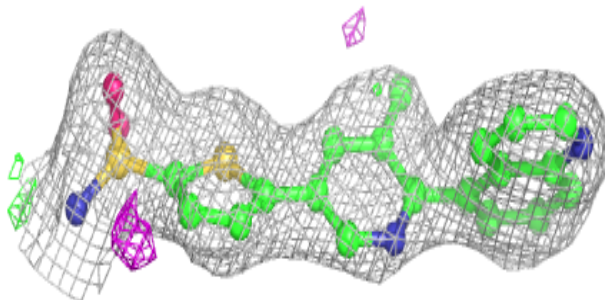
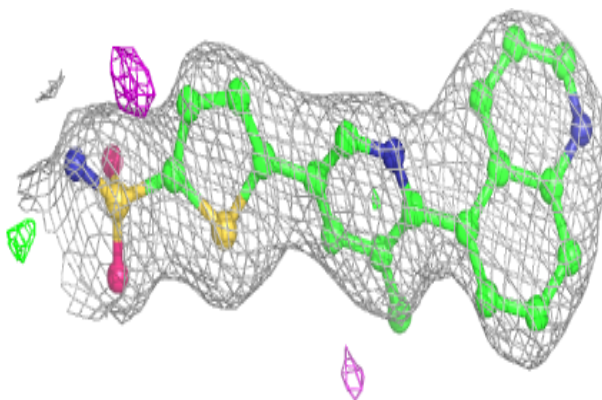


**Electron density around V80 C 302:**

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

**Electron density around V80 G 302:**

$2mF_o-DF_c$  (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.