



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 03:22 am BST

PDB ID : 6WGY  
Title : Crystal structure of a Putative citrate synthase 2 from *Mycobacterium bovis* in complex with citrate  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2020-04-06  
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 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

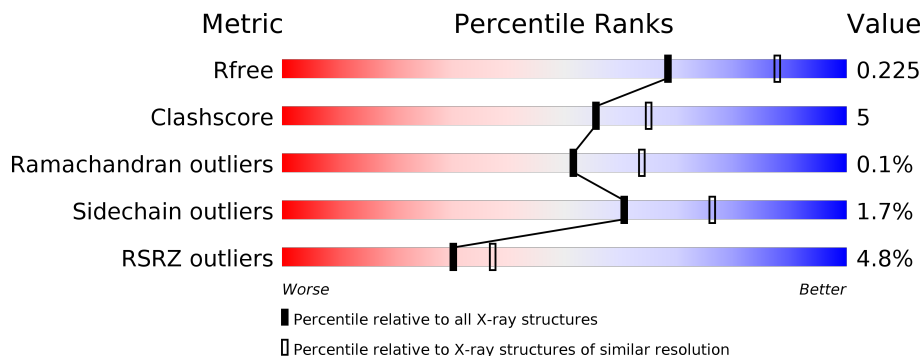
# 1 Overall quality at a glance

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 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	380	<p>7% 87% 12%</p>
1	B	380	<p>6% 89% 9%</p>
1	C	380	<p>5% 88% 9%</p>
1	D	380	<p>% 90% 8%</p>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	A	405	-	-	X	-
2	IOD	A	417	-	-	X	-
2	IOD	B	410	-	-	X	-
2	IOD	C	404	-	-	X	-
2	IOD	C	407	-	-	X	-
2	IOD	C	415	-	-	X	-
2	IOD	C	418	-	-	X	-
2	IOD	D	407[A]	-	-	X	-
2	IOD	D	413	-	-	X	-
2	IOD	D	414	-	-	X	-
2	IOD	D	416	-	-	X	-

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 12071 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 Putative citrate synthase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	374	2764	1748	490	512	14	0	2	0
1	B	374	2783	1759	497	513	14	0	2	0
1	C	371	2773	1753	492	513	15	0	4	0
1	D	373	2794	1764	495	521	14	0	2	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	374	GLY	-	expression tag	UNP P63778
A	375	HIS	-	expression tag	UNP P63778
A	376	HIS	-	expression tag	UNP P63778
A	377	HIS	-	expression tag	UNP P63778
A	378	HIS	-	expression tag	UNP P63778
A	379	HIS	-	expression tag	UNP P63778
A	380	HIS	-	expression tag	UNP P63778
B	374	GLY	-	expression tag	UNP P63778
B	375	HIS	-	expression tag	UNP P63778
B	376	HIS	-	expression tag	UNP P63778
B	377	HIS	-	expression tag	UNP P63778
B	378	HIS	-	expression tag	UNP P63778
B	379	HIS	-	expression tag	UNP P63778
B	380	HIS	-	expression tag	UNP P63778
C	374	GLY	-	expression tag	UNP P63778
C	375	HIS	-	expression tag	UNP P63778
C	376	HIS	-	expression tag	UNP P63778
C	377	HIS	-	expression tag	UNP P63778
C	378	HIS	-	expression tag	UNP P63778
C	379	HIS	-	expression tag	UNP P63778
C	380	HIS	-	expression tag	UNP P63778

*Continued on next page...*

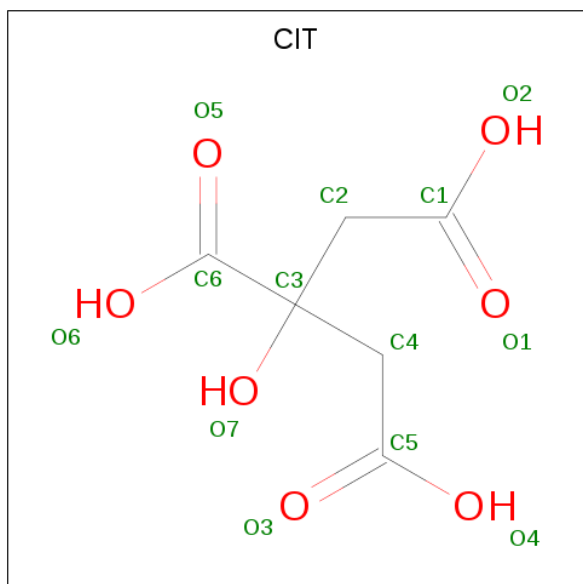
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	374	GLY	-	expression tag	UNP P63778
D	375	HIS	-	expression tag	UNP P63778
D	376	HIS	-	expression tag	UNP P63778
D	377	HIS	-	expression tag	UNP P63778
D	378	HIS	-	expression tag	UNP P63778
D	379	HIS	-	expression tag	UNP P63778
D	380	HIS	-	expression tag	UNP P63778

- Molecule 2 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	26	Total I 26 26	0	0
2	A	30	Total I 31 31	0	1
2	D	25	Total I 26 26	0	1
2	C	28	Total I 29 29	0	1

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: C<sub>6</sub>H<sub>8</sub>O<sub>7</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 13 6 7	0	0

Continued on next page...

*Continued from previous page...*

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			13	6	7		

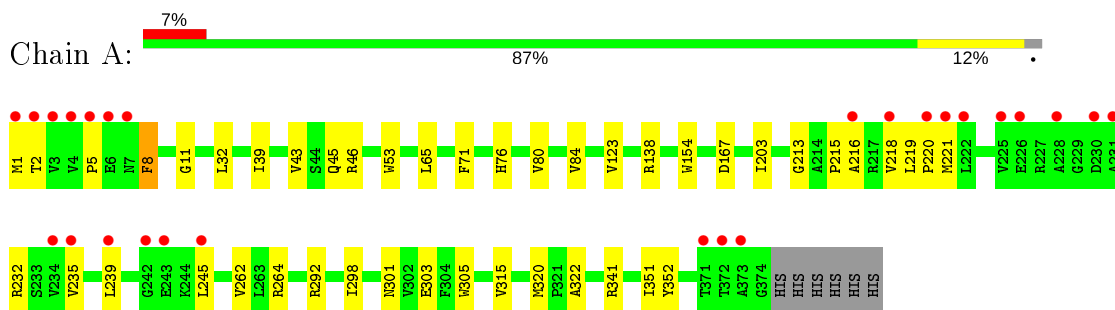
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	206	Total	O	0	6
			212	212		
4	B	180	Total	O	0	2
			182	182		
4	C	168	Total	O	0	3
			171	171		
4	D	248	Total	O	0	6
			254	254		

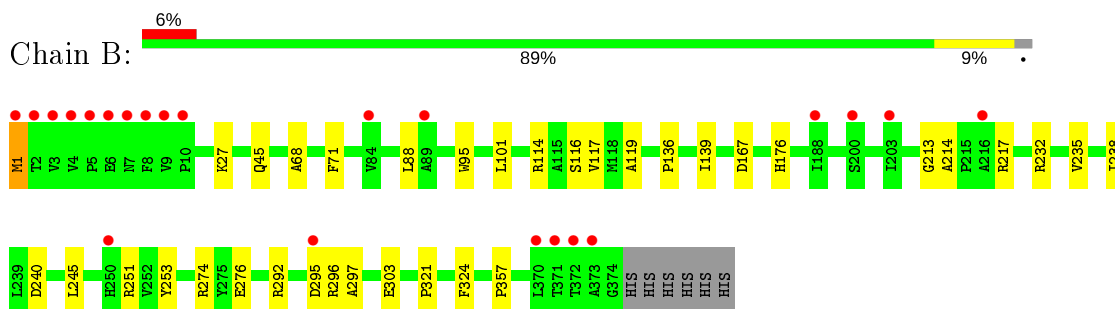
### 3 Residue-property plots [i](#)

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.

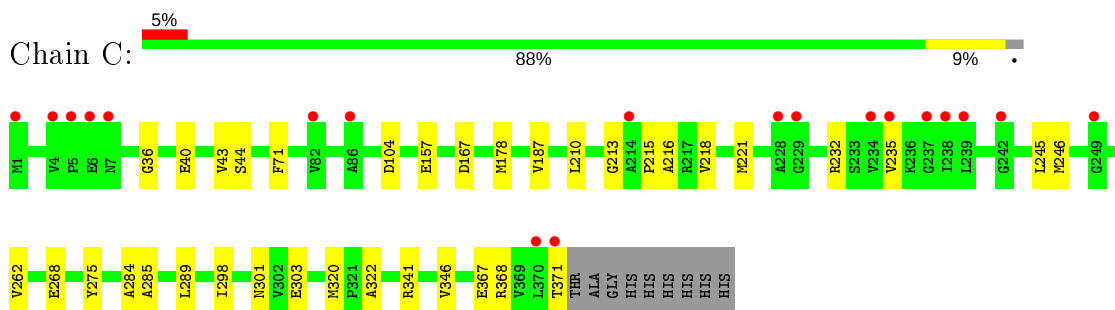
- Molecule 1: Putative citrate synthase 2



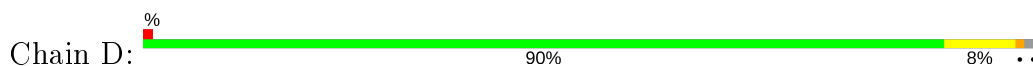
- Molecule 1: Putative citrate synthase 2

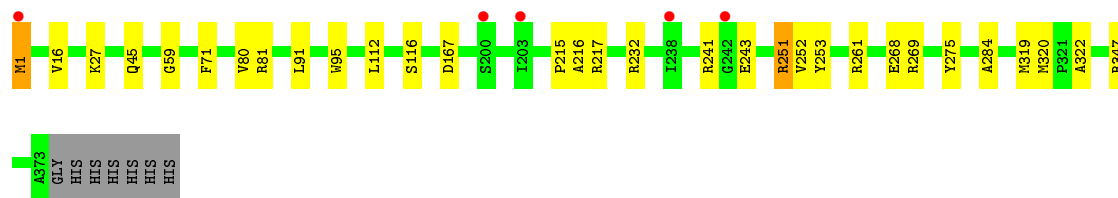


- Molecule 1: Putative citrate synthase 2



- Molecule 1: Putative citrate synthase 2







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.14Å 153.37Å 100.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.43 – 2.30 47.43 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.43-2.30) 100.0 (47.43-2.30)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.61 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.18rc4	Depositor
R, $R_{free}$	0.174 , 0.225 0.174 , 0.225	Depositor DCC
$R_{free}$ test set	1950 reflections (2.22%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.9	Xtrriage
Anisotropy	0.369	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12071	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: IOD, CIT

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	0.38	0/2834	0.53	0/3874
1	B	0.35	0/2854	0.52	0/3897
1	C	0.36	0/2846	0.52	0/3887
1	D	0.40	0/2864	0.53	0/3910
All	All	0.37	0/11398	0.52	0/15568

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2764	0	2694	32	0
1	B	2783	0	2724	24	0
1	C	2773	0	2707	30	0
1	D	2794	0	2741	23	0
2	A	31	0	0	14	0
2	B	26	0	0	7	0
2	C	29	0	0	15	0
2	D	26	0	0	9	0
3	B	13	0	5	1	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	13	0	5	0	0
4	A	212	0	0	4	0
4	B	182	0	0	2	0
4	C	171	0	0	2	0
4	D	254	0	0	6	0
All	All	12071	0	10876	112	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 5.

All (112) 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)
2:A:405:IOD:I	2:A:406:IOD:I	3.18	1.02
1:D:217:ARG:HA	2:D:414:IOD:I	2.47	0.85
1:B:251:ARG:HD2	1:B:297:ALA:HB3	1.61	0.82
1:D:322:ALA:HA	2:D:407[A]:IOD:I	2.62	0.70
1:B:213:GLY:O	2:B:411:IOD:I	2.81	0.68
1:A:322:ALA:HA	2:A:405:IOD:I	2.63	0.67
2:A:427:IOD:I	2:C:424:IOD:I	3.54	0.66
1:B:213:GLY:O	1:B:303:GLU:HG3	1.96	0.65
1:A:213:GLY:O	2:A:415:IOD:I	2.87	0.63
1:A:235:VAL:HG13	1:A:245:LEU:HD11	1.81	0.62
1:B:235:VAL:HG13	1:B:245:LEU:HD11	1.80	0.62
1:D:81:ARG:HA	2:D:407[A]:IOD:I	2.69	0.62
2:A:426:IOD:I	2:C:427:IOD:I	3.56	0.62
1:A:220:PRO:HD3	2:A:417:IOD:I	2.70	0.62
1:D:261:ARG:NH1	4:D:504:HOH:O	2.32	0.61
1:C:322:ALA:HA	2:C:407:IOD:I	2.70	0.61
1:D:80:VAL:HG21	4:D:734:HOH:O	2.00	0.60
2:A:412:IOD:I	2:D:401:IOD:I	3.60	0.60
1:B:214:ALA:HB1	2:B:410:IOD:I	2.72	0.59
1:A:341:ARG:NH1	4:A:505:HOH:O	2.33	0.59
1:B:321:PRO:HG3	2:B:410:IOD:I	2.72	0.58
1:C:289:LEU:HB3	1:C:298:ILE:HD13	1.84	0.58
1:B:176:HIS:ND1	3:B:426:CIT:H42	2.18	0.58
1:A:45:GLN:NE2	4:A:504:HOH:O	2.32	0.57
1:C:213:GLY:O	2:C:415:IOD:I	2.92	0.57
1:A:1:MET:HG2	1:A:352:TYR:H	1.68	0.57
1:D:45:GLN:NE2	4:D:506:HOH:O	2.35	0.56
1:A:43:VAL:HG11	1:A:262:VAL:HG13	1.88	0.56

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:ALA:O	2:D:414:IOD:I	2.95	0.55
1:C:341:ARG:NH1	4:C:501:HOH:O	2.27	0.55
1:B:1:MET:HG3	1:B:1:MET:O	2.07	0.54
1:A:232:ARG:HD3	2:A:419:IOD:I	2.78	0.54
1:A:46:ARG:HD3	4:D:502:HOH:O	2.07	0.54
1:C:215:PRO:HG2	1:C:320:MET:HB3	1.90	0.53
1:A:219:LEU:HB2	2:A:417:IOD:I	2.78	0.53
1:A:84[A]:VAL:HG12	1:A:203:ILE:HD13	1.91	0.53
1:C:232:ARG:HD3	2:C:418:IOD:I	2.78	0.53
2:C:407:IOD:I	4:C:651:HOH:O	2.89	0.52
1:C:43:VAL:HG11	1:C:262:VAL:HG13	1.91	0.52
1:C:235:VAL:HG11	1:C:285:ALA:HA	1.92	0.52
1:A:11:GLY:O	4:A:501:HOH:O	2.19	0.51
1:D:216:ALA:HB2	2:D:413:IOD:I	2.81	0.51
1:C:303:GLU:HG3	2:C:415:IOD:I	2.81	0.51
1:C:104:ASP:HB2	2:C:428:IOD:I	2.81	0.51
1:A:351:ILE:HG12	1:D:16:VAL:HG11	1.93	0.50
1:C:44:SER:HB3	2:C:404:IOD:I	2.81	0.50
1:C:235:VAL:HG13	1:C:245:LEU:HD11	1.93	0.50
1:A:262:VAL:HG22	2:A:402:IOD:I	2.82	0.50
1:A:215:PRO:HG2	1:A:320:MET:HB3	1.93	0.50
1:D:112:LEU:O	1:D:116:SER:HB2	2.11	0.50
1:D:269:ARG:HD3	2:D:416:IOD:I	2.82	0.49
1:C:40:GLU:HG2	1:C:262:VAL:HG23	1.95	0.49
1:B:45:GLN:O	1:C:371:THR:OG1	2.31	0.49
1:B:27:LYS:HA	1:B:253:TYR:CD1	2.48	0.49
1:D:1:MET:HG3	1:D:1:MET:O	2.13	0.49
1:A:239:LEU:HB3	1:A:292:ARG:HH12	1.77	0.48
1:C:268:GLU:HB2	1:C:275:TYR:CZ	2.48	0.48
1:D:241:ARG:NE	1:D:243:GLU:OE1	2.40	0.47
1:C:268:GLU:HB2	1:C:275:TYR:CE1	2.49	0.47
1:B:238:ILE:HD11	2:B:414:IOD:I	2.85	0.47
1:A:215:PRO:O	2:A:414:IOD:I	3.03	0.47
1:A:76:HIS:ND1	4:A:506:HOH:O	2.35	0.46
1:B:136:PRO:HD2	1:B:139[A]:ILE:HD12	1.98	0.46
1:C:215:PRO:O	2:C:414:IOD:I	3.04	0.46
1:C:221:MET:HA	2:C:417:IOD:I	2.86	0.46
2:C:404:IOD:I	2:C:419:IOD:I	3.74	0.46
1:D:216:ALA:CB	2:D:413:IOD:I	3.34	0.45
1:A:5:PRO:HG2	1:A:8:PHE:HB2	1.97	0.45
1:B:95:TRP:CZ3	1:B:114:ARG:HG2	2.51	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:367:GLU:HG2	1:C:368:ARG:N	2.32	0.45
1:B:274:ARG:HG2	2:B:423:IOD:I	2.87	0.45
1:B:68:ALA:HB2	1:B:117:VAL:HG22	1.99	0.44
1:C:216:ALA:CB	1:C:218:VAL:HG13	2.48	0.44
1:C:301:ASN:HB3	1:C:303:GLU:OE1	2.18	0.44
1:A:32:LEU:HD23	1:A:39:ILE:HD13	2.00	0.44
1:D:268:GLU:HB2	1:D:275:TYR:CZ	2.53	0.44
1:A:154:TRP:HH2	1:A:315:VAL:HA	1.83	0.43
1:B:296:ARG:NH2	2:B:424:IOD:I	3.22	0.43
1:C:218:VAL:HG12	1:C:246:MET:HB2	2.00	0.43
1:D:215:PRO:HG2	1:D:320:MET:HB3	2.01	0.43
2:D:404:IOD:I	2:D:416:IOD:I	3.76	0.43
1:A:232:ARG:HD2	1:A:232:ARG:HA	1.73	0.43
1:B:1:MET:N	4:B:511:HOH:O	2.52	0.43
1:C:216:ALA:HB2	2:C:416:IOD:I	2.89	0.43
1:A:84[A]:VAL:HG21	1:A:123:VAL:HG22	2.01	0.42
1:C:216:ALA:HB1	1:C:218:VAL:HG13	2.01	0.42
1:B:357:PRO:HA	1:C:36:GLY:O	2.18	0.42
1:B:295:ASP:HB2	2:B:417:IOD:I	2.89	0.42
1:B:101:LEU:HD13	1:C:210:LEU:HD11	2.02	0.42
1:D:27:LYS:HD2	1:D:252:VAL:O	2.18	0.42
1:A:301:ASN:HB3	1:A:303:GLU:OE1	2.20	0.42
1:B:213:GLY:HA3	1:B:324:PHE:CE2	2.54	0.42
1:D:91:LEU:HD13	1:D:95:TRP:CZ3	2.55	0.42
1:D:27:LYS:HA	1:D:253:TYR:CD1	2.55	0.42
1:A:216:ALA:HB1	1:A:218:VAL:HG13	2.02	0.42
1:A:298:ILE:HG12	2:A:429[A]:IOD:I	2.90	0.42
1:A:80[A]:VAL:HG21	2:A:423:IOD:I	2.90	0.41
1:A:221:MET:HA	2:A:418:IOD:I	2.90	0.41
1:A:53:TRP:CD2	1:A:65:LEU:HD13	2.54	0.41
1:D:59:GLY:O	4:D:501:HOH:O	2.22	0.41
1:C:187:VAL:HG22	1:C:346:VAL:HG11	2.01	0.41
1:D:251:ARG:HB2	1:D:253:TYR:O	2.21	0.41
1:A:264:ARG:HG3	1:A:305:TRP:CZ2	2.55	0.41
1:B:27:LYS:NZ	4:B:512:HOH:O	2.53	0.41
1:A:84[A]:VAL:HG21	1:A:123:VAL:CG2	2.50	0.41
1:D:319:MET:HG3	4:D:734:HOH:O	2.20	0.41
1:B:240:ASP:OD1	1:B:292:ARG:NH2	2.48	0.41
1:C:284:ALA:HA	2:C:418:IOD:I	2.91	0.41
1:C:341:ARG:NH2	2:C:409:IOD:I	3.24	0.41
1:D:232:ARG:HA	1:D:284:ALA:HB1	2.02	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:232:ARG:HA	1:C:284:ALA:HB1	2.03	0.40
1:B:88:LEU:HD11	1:B:119:ALA:HB2	2.02	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	374/380 (98%)	363 (97%)	10 (3%)	1 (0%)	41	50
1	B	374/380 (98%)	359 (96%)	14 (4%)	1 (0%)	41	50
1	C	373/380 (98%)	364 (98%)	9 (2%)	0	100	100
1	D	373/380 (98%)	364 (98%)	9 (2%)	0	100	100
All	All	1494/1520 (98%)	1450 (97%)	42 (3%)	2 (0%)	51	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	217	ARG
1	A	2	THR

### 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	271/291 (93%)	267 (98%)	4 (2%)	65	79
1	B	274/291 (94%)	268 (98%)	6 (2%)	52	69
1	C	274/291 (94%)	268 (98%)	6 (2%)	52	69
1	D	279/291 (96%)	274 (98%)	5 (2%)	59	75
All	All	1098/1164 (94%)	1077 (98%)	21 (2%)	60	73

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	71	PHE
1	A	138	ARG
1	A	167	ASP
1	B	1	MET
1	B	71	PHE
1	B	116	SER
1	B	167	ASP
1	B	232	ARG
1	B	276	GLU
1	C	71	PHE
1	C	157	GLU
1	C	167[A]	ASP
1	C	167[B]	ASP
1	C	178[A]	MET
1	C	178[B]	MET
1	D	1	MET
1	D	71	PHE
1	D	167	ASP
1	D	251	ARG
1	D	347	ARG

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

Mol	Chain	Res	Type
1	A	76	HIS
1	B	179	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 114 ligands modelled in this entry, 112 are monoatomic - leaving 2 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	CIT	B	426	-	3,12,12	1.55	0	3,17,17	2.08	1 (33%)
3	CIT	D	426	-	3,12,12	1.61	0	3,17,17	2.27	2 (66%)

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	CIT	B	426	-	-	2/6/16/16	-
3	CIT	D	426	-	-	0/6/16/16	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	426	CIT	C3-C4-C5	-3.55	109.31	114.98
3	D	426	CIT	C3-C4-C5	2.77	119.42	114.98
3	D	426	CIT	C3-C2-C1	-2.30	111.31	114.98

There are no chirality outliers.



All (2) torsion outliers are listed below:

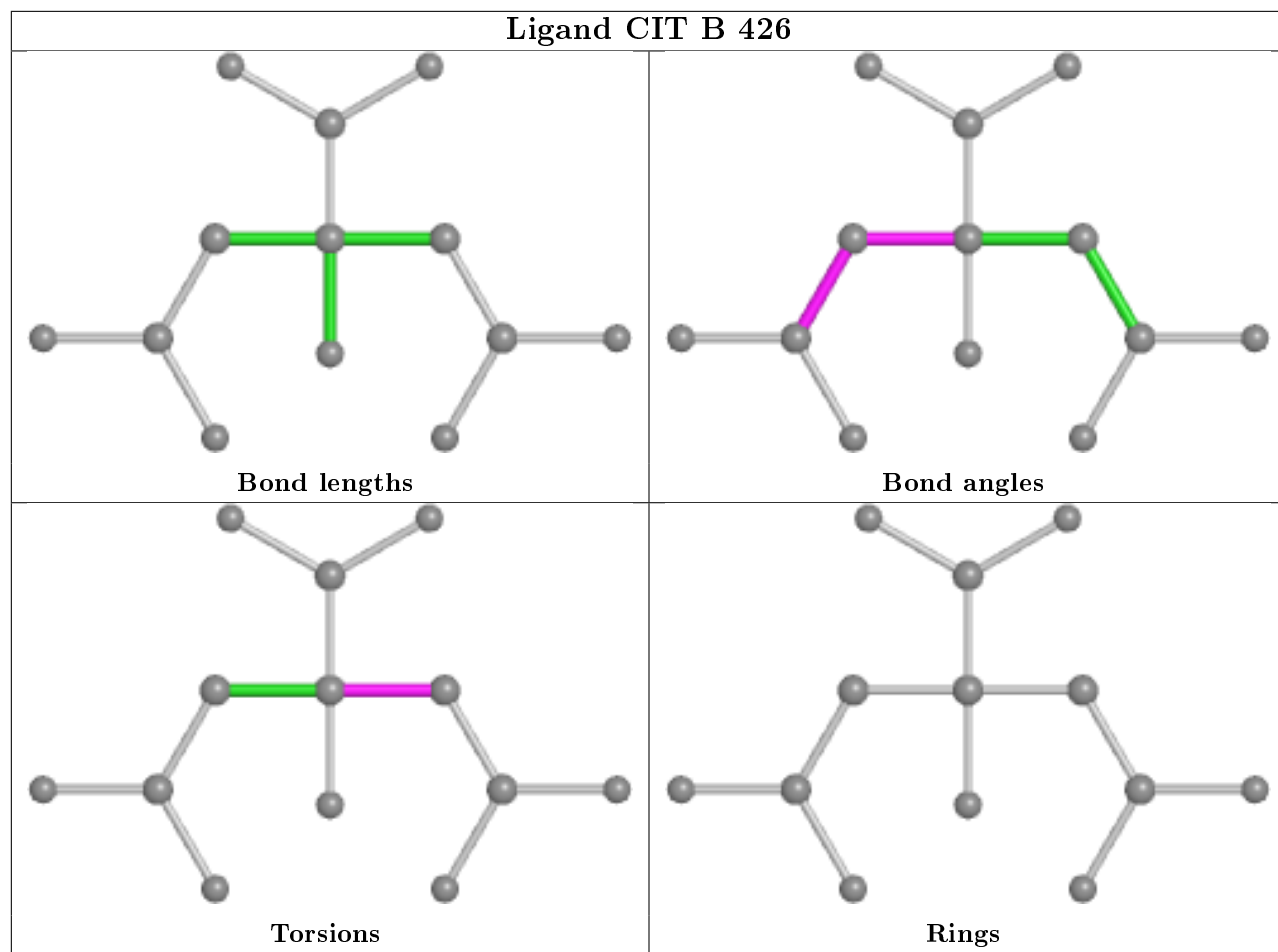
Mol	Chain	Res	Type	Atoms
3	B	426	CIT	C1-C2-C3-C4
3	B	426	CIT	C1-C2-C3-C6

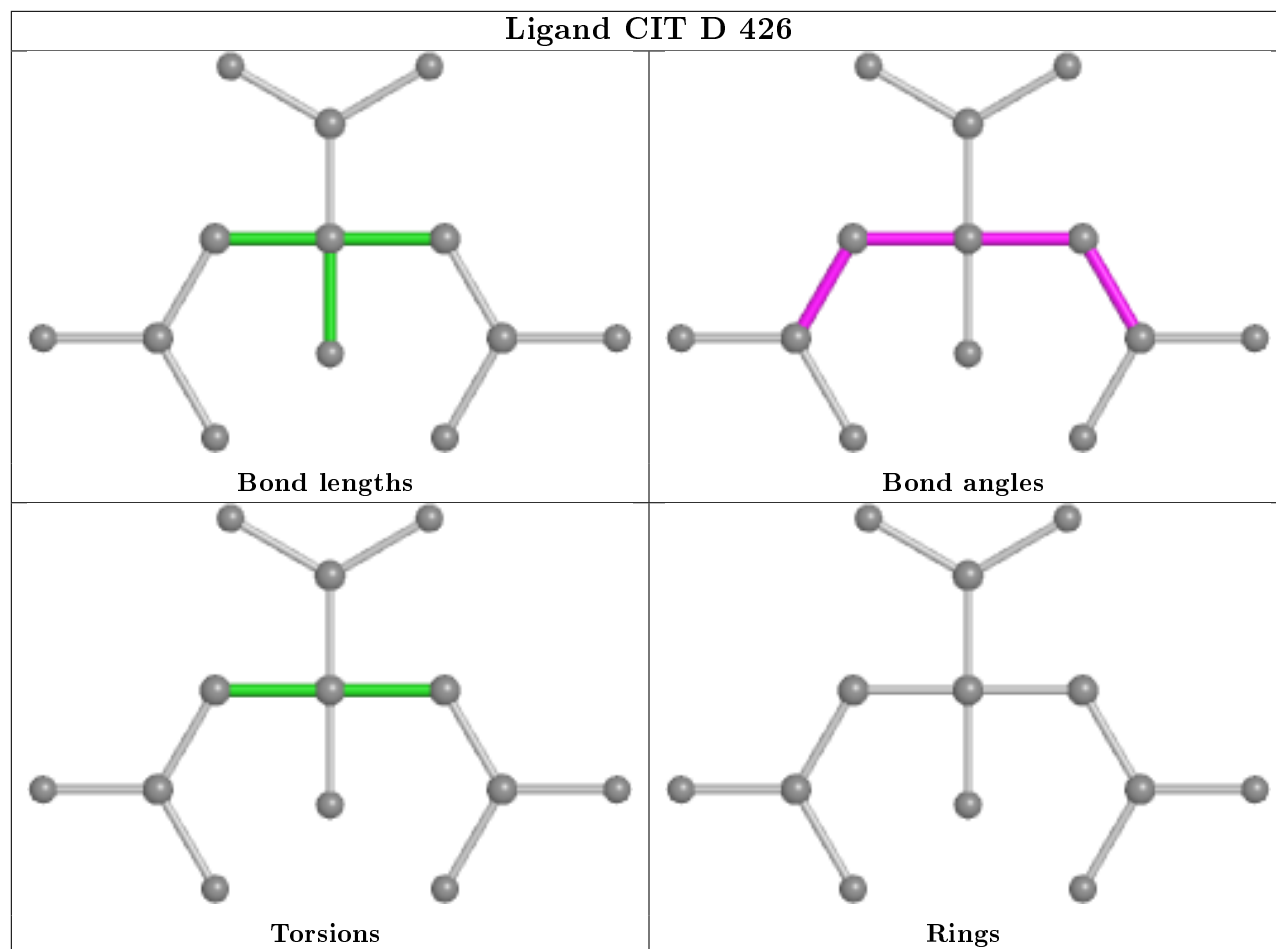
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	426	CIT	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 [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	374/380 (98%)	0.07	26 (6%) 16 21	23, 38, 87, 112	0
1	B	374/380 (98%)	0.06	22 (5%) 22 28	25, 40, 74, 106	0
1	C	371/380 (97%)	0.03	19 (5%) 28 35	26, 46, 86, 113	0
1	D	373/380 (98%)	-0.11	5 (1%) 77 81	23, 35, 62, 96	0
All	All	1492/1520 (98%)	0.01	72 (4%) 30 37	23, 39, 81, 113	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	371	THR	6.8
1	A	235	VAL	6.7
1	C	242	GLY	5.1
1	B	4	VAL	5.0
1	B	8	PHE	4.9
1	B	216	ALA	4.5
1	C	4	VAL	4.3
1	A	1	MET	4.1
1	C	1	MET	4.1
1	C	239	LEU	4.1
1	B	10	PRO	3.9
1	A	239	LEU	3.9
1	A	7	ASN	3.7
1	B	373	ALA	3.7
1	B	372	THR	3.7
1	A	234	VAL	3.6
1	A	225	VAL	3.6
1	C	5	PRO	3.4
1	A	242	GLY	3.3
1	A	230	ASP	3.3
1	D	1	MET	3.2

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	234	VAL	3.1
1	C	371	THR	3.1
1	C	7	ASN	3.1
1	B	5	PRO	3.0
1	D	242	GLY	2.9
1	C	229	GLY	2.9
1	B	7	ASN	2.9
1	A	231	ALA	2.8
1	B	6	GLU	2.7
1	C	235	VAL	2.7
1	A	216	ALA	2.6
1	A	4	VAL	2.6
1	A	218	VAL	2.6
1	A	2	THR	2.5
1	B	1	MET	2.5
1	D	203	ILE	2.5
1	A	226	GLU	2.5
1	C	82	VAL	2.5
1	C	237	GLY	2.5
1	D	238	ILE	2.5
1	B	89	ALA	2.4
1	B	188	ILE	2.4
1	A	373	ALA	2.4
1	C	370	LEU	2.4
1	B	250[A]	HIS	2.4
1	C	238	ILE	2.3
1	A	372	THR	2.3
1	C	214	ALA	2.3
1	A	220	PRO	2.3
1	A	5	PRO	2.3
1	D	200	SER	2.3
1	B	3	VAL	2.2
1	A	371	THR	2.2
1	A	6	GLU	2.1
1	A	3	VAL	2.1
1	C	228	ALA	2.1
1	A	222	LEU	2.1
1	A	245	LEU	2.1
1	A	243	GLU	2.1
1	C	6	GLU	2.1
1	B	2	THR	2.1
1	B	9	VAL	2.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	370	LEU	2.1
1	A	228	ALA	2.1
1	B	203	ILE	2.1
1	B	295	ASP	2.1
1	B	200	SER	2.1
1	A	221	MET	2.0
1	B	84	VAL	2.0
1	C	249	GLY	2.0
1	C	86	ALA	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 carbohydrates 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
2	IOD	A	414	1/1	0.76	0.10	109,109,109,109	1
2	IOD	B	425	1/1	0.77	0.26	210,210,210,210	0
2	IOD	A	406	1/1	0.78	0.19	68,68,68,68	1
2	IOD	A	416	1/1	0.78	0.11	117,117,117,117	1
2	IOD	A	417	1/1	0.79	0.18	109,109,109,109	1
2	IOD	C	424	1/1	0.80	0.13	91,91,91,91	1
2	IOD	B	420	1/1	0.83	0.10	86,86,86,86	1
3	CIT	B	426	13/13	0.84	0.19	44,57,66,67	0
2	IOD	C	416	1/1	0.85	0.08	97,97,97,97	1
2	IOD	C	426[A]	1/1	0.86	0.09	104,104,104,104	1
2	IOD	A	429[A]	1/1	0.86	0.08	113,113,113,113	1
2	IOD	A	429[B]	1/1	0.86	0.08	113,113,113,113	1
2	IOD	C	426[B]	1/1	0.86	0.09	99,99,99,99	1
2	IOD	D	424	1/1	0.87	0.05	90,90,90,90	1

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	IOD	B	416	1/1	0.87	0.08	104,104,104,104	1
2	IOD	A	415	1/1	0.88	0.10	90,90,90,90	1
2	IOD	C	401	1/1	0.88	0.11	55,55,55,55	1
2	IOD	A	430	1/1	0.88	0.11	165,165,165,165	1
2	IOD	C	403	1/1	0.89	0.08	83,83,83,83	1
2	IOD	D	422	1/1	0.89	0.21	89,89,89,89	1
3	CIT	D	426	13/13	0.90	0.16	34,44,50,53	0
2	IOD	C	402	1/1	0.90	0.06	71,71,71,71	0
2	IOD	A	428	1/1	0.90	0.10	89,89,89,89	1
2	IOD	A	418	1/1	0.90	0.06	86,86,86,86	1
2	IOD	B	423	1/1	0.91	0.07	79,79,79,79	1
2	IOD	D	418	1/1	0.91	0.06	77,77,77,77	1
2	IOD	D	416	1/1	0.92	0.07	85,85,85,85	1
2	IOD	B	415	1/1	0.92	0.15	99,99,99,99	1
2	IOD	C	414	1/1	0.92	0.10	97,97,97,97	1
2	IOD	A	421	1/1	0.92	0.11	115,115,115,115	1
2	IOD	C	428	1/1	0.92	0.07	118,118,118,118	1
2	IOD	B	401	1/1	0.92	0.05	92,92,92,92	1
2	IOD	A	420	1/1	0.92	0.08	89,89,89,89	1
2	IOD	B	421	1/1	0.92	0.09	95,95,95,95	1
2	IOD	B	417	1/1	0.93	0.15	111,111,111,111	1
2	IOD	C	406	1/1	0.93	0.08	57,57,57,57	1
2	IOD	A	426	1/1	0.93	0.13	68,68,68,68	1
2	IOD	D	406	1/1	0.93	0.12	66,66,66,66	1
2	IOD	B	409	1/1	0.93	0.07	90,90,90,90	1
2	IOD	D	412	1/1	0.93	0.12	107,107,107,107	1
2	IOD	A	412	1/1	0.94	0.15	105,105,105,105	1
2	IOD	D	415	1/1	0.94	0.04	72,72,72,72	1
2	IOD	B	419	1/1	0.94	0.07	67,67,67,67	1
2	IOD	C	415	1/1	0.94	0.13	84,84,84,84	1
2	IOD	D	414	1/1	0.94	0.36	105,105,105,105	1
2	IOD	D	403	1/1	0.94	0.06	78,78,78,78	1
2	IOD	C	425	1/1	0.94	0.10	89,89,89,89	1
2	IOD	B	411	1/1	0.94	0.11	75,75,75,75	1
2	IOD	A	404	1/1	0.95	0.11	65,65,65,65	1
2	IOD	B	412	1/1	0.95	0.05	77,77,77,77	1
2	IOD	D	425	1/1	0.95	0.07	95,95,95,95	1
2	IOD	B	404	1/1	0.95	0.09	55,55,55,55	1
2	IOD	D	420	1/1	0.95	0.12	78,78,78,78	1
2	IOD	C	407	1/1	0.95	0.11	46,46,46,46	1
2	IOD	B	410	1/1	0.96	0.07	97,97,97,97	1
2	IOD	D	419	1/1	0.96	0.09	61,61,61,61	1

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	IOD	B	427	1/1	0.96	0.08	96,96,96,96	1
2	IOD	D	421	1/1	0.96	0.04	69,69,69,69	1
2	IOD	A	411	1/1	0.96	0.04	74,74,74,74	1
2	IOD	B	405	1/1	0.97	0.06	70,70,70,70	1
2	IOD	D	407[A]	1/1	0.97	0.07	60,60,60,60	1
2	IOD	A	401	1/1	0.97	0.04	52,52,52,52	1
2	IOD	D	411	1/1	0.97	0.04	74,74,74,74	1
2	IOD	A	410	1/1	0.97	0.07	62,62,62,62	1
2	IOD	B	418	1/1	0.97	0.07	60,60,60,60	1
2	IOD	C	408	1/1	0.97	0.08	68,68,68,68	1
2	IOD	B	414	1/1	0.97	0.06	80,80,80,80	1
2	IOD	A	423	1/1	0.97	0.06	74,74,74,74	1
2	IOD	C	417	1/1	0.97	0.04	85,85,85,85	1
2	IOD	C	423	1/1	0.97	0.06	65,65,65,65	1
2	IOD	A	405	1/1	0.97	0.13	44,44,44,44	1
2	IOD	D	407[B]	1/1	0.97	0.07	62,62,62,62	1
2	IOD	A	422	1/1	0.97	0.06	80,80,80,80	1
2	IOD	A	427	1/1	0.97	0.12	89,89,89,89	1
2	IOD	C	411	1/1	0.97	0.04	65,65,65,65	1
2	IOD	C	412	1/1	0.97	0.08	90,90,90,90	1
2	IOD	B	402	1/1	0.97	0.07	54,54,54,54	1
2	IOD	D	401	1/1	0.98	0.09	41,41,41,41	1
2	IOD	C	413	1/1	0.98	0.08	59,59,59,59	1
2	IOD	B	406	1/1	0.98	0.06	55,55,55,55	1
2	IOD	C	420	1/1	0.98	0.06	94,94,94,94	1
2	IOD	C	410	1/1	0.98	0.10	53,53,53,53	1
2	IOD	D	413	1/1	0.98	0.07	61,61,61,61	1
2	IOD	B	413	1/1	0.98	0.05	76,76,76,76	1
2	IOD	D	404	1/1	0.98	0.06	52,52,52,52	1
2	IOD	C	409	1/1	0.98	0.09	72,72,72,72	1
2	IOD	B	422	1/1	0.98	0.11	89,89,89,89	1
2	IOD	A	419	1/1	0.98	0.05	96,96,96,96	1
2	IOD	C	419	1/1	0.98	0.06	94,94,94,94	1
2	IOD	D	408	1/1	0.98	0.07	56,56,56,56	1
2	IOD	C	421	1/1	0.98	0.03	74,74,74,74	1
2	IOD	B	424	1/1	0.98	0.04	68,68,68,68	1
2	IOD	A	424	1/1	0.98	0.11	46,46,46,46	1
2	IOD	D	402	1/1	0.99	0.09	97,97,97,97	1
2	IOD	D	405	1/1	0.99	0.09	47,47,47,47	1
2	IOD	A	409	1/1	0.99	0.08	43,43,43,43	1
2	IOD	D	409	1/1	0.99	0.09	51,51,51,51	1
2	IOD	A	425	1/1	0.99	0.04	64,64,64,64	1

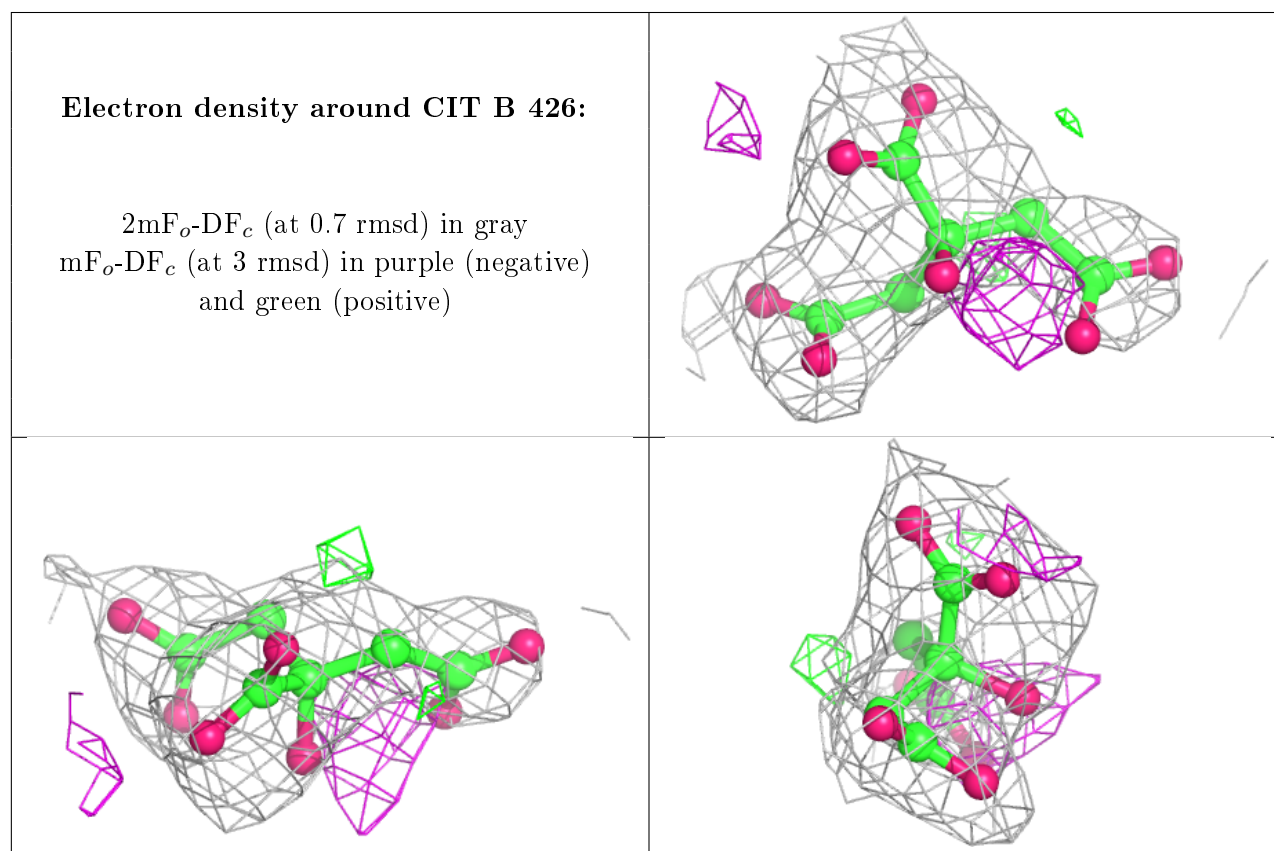
*Continued on next page...*

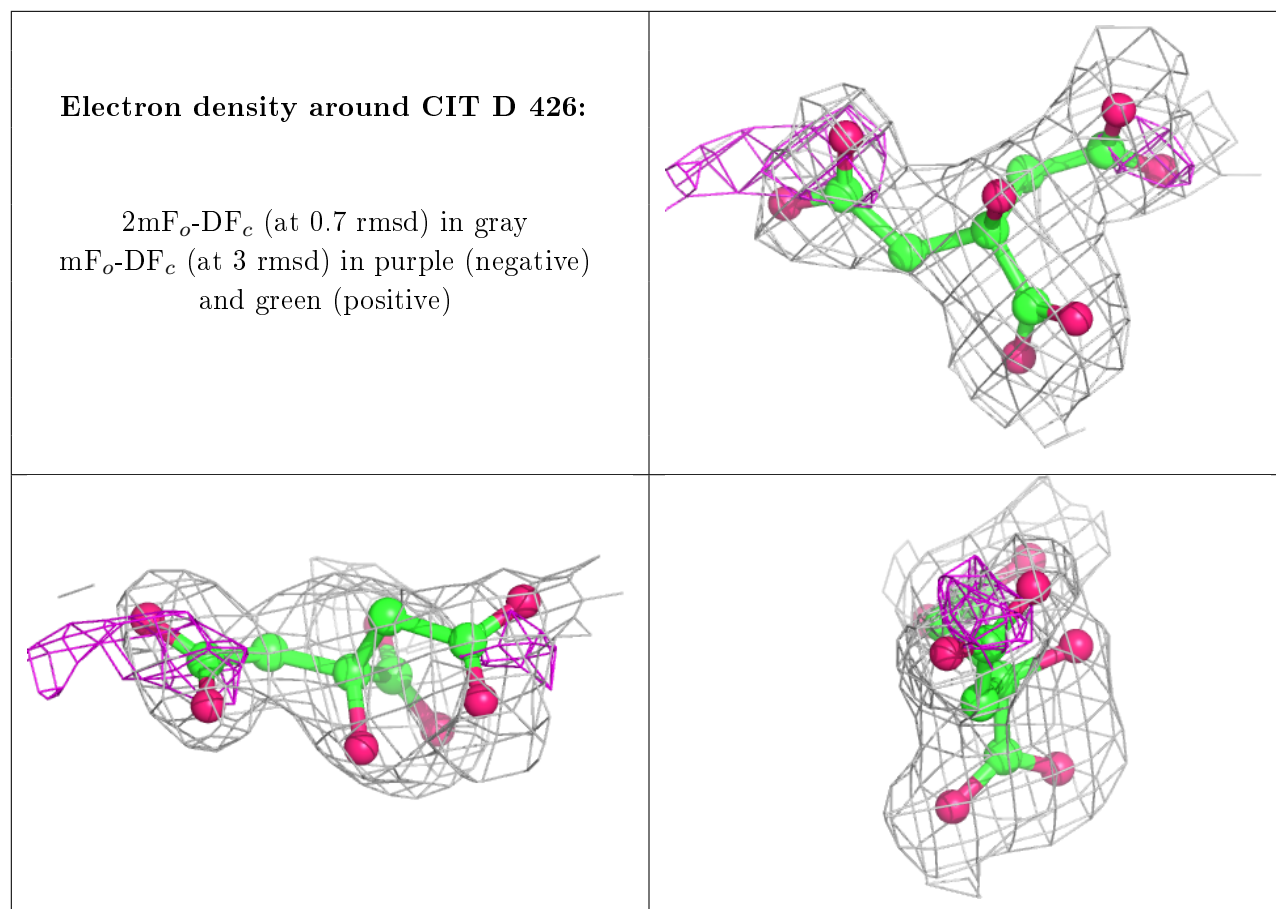


Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	IOD	B	407	1/1	0.99	0.09	49,49,49,49	1
2	IOD	C	404	1/1	0.99	0.04	67,67,67,67	1
2	IOD	A	402	1/1	0.99	0.06	61,61,61,61	1
2	IOD	A	408	1/1	0.99	0.06	46,46,46,46	1
2	IOD	D	417	1/1	0.99	0.09	86,86,86,86	1
2	IOD	B	403	1/1	0.99	0.06	53,53,53,53	1
2	IOD	D	410	1/1	0.99	0.08	42,42,42,42	1
2	IOD	C	418	1/1	0.99	0.04	75,75,75,75	1
2	IOD	C	422	1/1	0.99	0.05	54,54,54,54	1
2	IOD	C	427	1/1	0.99	0.10	77,77,77,77	1
2	IOD	C	405	1/1	0.99	0.05	55,55,55,55	0
2	IOD	B	408	1/1	0.99	0.04	60,60,60,60	1
2	IOD	A	407	1/1	0.99	0.07	48,48,48,48	1
2	IOD	D	423	1/1	0.99	0.04	73,73,73,73	1
2	IOD	A	413	1/1	0.99	0.11	50,50,50,50	1
2	IOD	A	403	1/1	1.00	0.08	46,46,46,46	1

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.





## 6.5 Other polymers [i](#)

There are no such residues in this entry.