

# Full wwPDB X-ray Structure Validation Report (i)

#### Feb 15, 2024 – 08:16 PM EST

PDB ID : 3R01

Title: The discovery of novel benzofuran-2-carboxylic acids as potent Pim-1 inhibitors

Authors : Liu, J. Deposited on : 2011-03-07

Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), 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)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

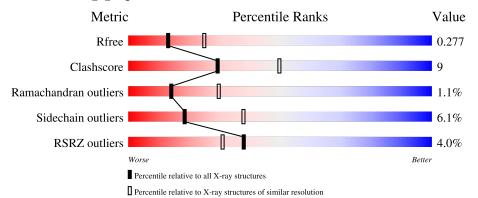
Validation Pipeline (wwPDB-VP) : 2.36

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY\ DIFFRACTION$ 

The reported resolution of this entry is 2.60 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	Similar resolution
Metric	$(\#  ext{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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			
			4%			
1	A	299	73%	13%	5%	8%



## 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2346 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 Proto-oncogene serine/threonine-protein kinase pim-1.

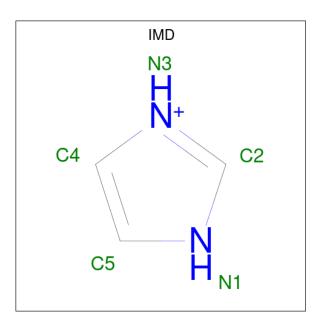
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	274	Total	С	N	О	S	0	0	0
1	A	214	2230	1426	390	406	8	0	U	

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	MET	-	initiating methionine	UNP P11309
A	314	ALA	-	expression tag	UNP P11309
A	315	ALA	-	expression tag	UNP P11309
A	316	ALA	-	expression tag	UNP P11309
A	317	LEU	-	expression tag	UNP P11309
A	318	GLU	-	expression tag	UNP P11309
A	319	HIS	-	expression tag	UNP P11309
A	320	HIS	-	expression tag	UNP P11309
A	321	HIS	-	expression tag	UNP P11309
A	322	HIS	-	expression tag	UNP P11309
A	323	HIS	-	expression tag	UNP P11309
A	324	HIS	-	expression tag	UNP P11309
A	325	HIS	-	expression tag	UNP P11309
A	326	HIS	-	expression tag	UNP P11309

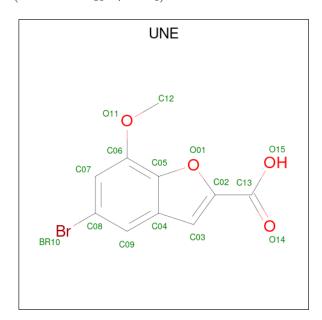
• Molecule 2 is IMIDAZOLE (three-letter code: IMD) (formula:  $C_3H_5N_2$ ).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N 5 3 2	0	0
2	A	1	Total C N 5 3 2	0	0

• Molecule 3 is 5-bromo-7-methoxy-1-benzofuran-2-carboxylic acid (three-letter code: UNE) (formula:  $C_{10}H_7BrO_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 15	Br 1	C 10	O 4	0	0



• Molecule 4 is water.

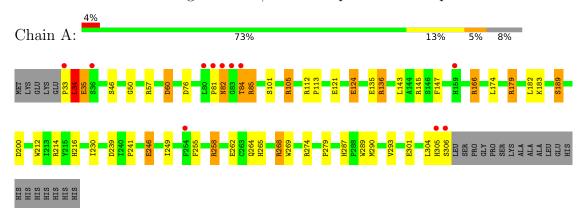
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	91	Total O 91 91	0	0



## 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: Proto-oncogene serine/threonine-protein kinase pim-1





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants	97.94Å 97.94Å 80.75Å	Donositon
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor
Resolution (Å)	84.82 - 2.60	Depositor
Resolution (A)	32.06 - 2.60	EDS
% Data completeness	99.9 (84.82-2.60)	Depositor
(in resolution range)	100.0 (32.06-2.60)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	4.83  (at  2.61Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
$R, R_{free}$	0.215 , $0.287$	Depositor
it, it free	0.210 , $0.277$	DCC
$R_{free}$ test set	677  reflections  (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.2	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.38, 43.6	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.54, < L^2> = 0.38$	Xtriage
Estimated twinning fraction	0.023 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2346	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.

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



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

## 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: UNE, IMD

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).

Mal	Chain	Bo	nd lengths	Bond angles		
Mol   Cha	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	1.02	$2/2290 \ (0.1\%)$	0.97	7/3108 (0.2%)	

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

#### All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}( ext{\AA})$
1	A	258	ARG	CZ-NH1	6.81	1.42	1.33
1	A	301	GLU	CG-CD	5.10	1.59	1.51

#### All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
1	A	268	ARG	NE-CZ-NH1	-10.62	114.99	120.30
1	A	268	ARG	NE-CZ-NH2	8.93	124.76	120.30
1	A	166	ARG	NE-CZ-NH2	-7.59	116.50	120.30
1	A	258	ARG	NE-CZ-NH1	7.05	123.82	120.30
1	A	34	LEU	CA-CB-CG	6.33	129.87	115.30
1	A	166	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	A	258	ARG	NE-CZ-NH2	-5.65	117.47	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:



Mol	Chain	Res	Type	Group	
1	A	35	GLU	Peptide	

#### 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	2230	0	2175	41	0
2	A	10	0	10	1	0
3	A	15	0	6	1	0
4	A	91	0	0	8	0
All	All	2346	0	2191	42	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 9.

All (42) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:268:ARG:HD2	4:A:25:HOH:O	1.56	1.03
1:A:265:HIS:HD2	1:A:287:HIS:HE1	0.95	0.93
1:A:265:HIS:HD2	1:A:287:HIS:CE1	1.87	0.92
1:A:287:HIS:HD2	1:A:289:TRP:H	1.13	0.92
1:A:265:HIS:CD2	1:A:287:HIS:HE1	1.87	0.91
1:A:262:GLU:HG3	4:A:357:HOH:O	1.69	0.90
1:A:105:ARG:HG2	1:A:121:GLU:OE1	1.76	0.86
1:A:124:GLU:HG2	4:A:8:HOH:O	1.86	0.75
1:A:274:ARG:HD3	4:A:343:HOH:O	1.87	0.73
1:A:287:HIS:CD2	1:A:289:TRP:H	2.03	0.72
1:A:166:ARG:HD3	1:A:189:SER:O	1.90	0.70
1:A:214:ARG:CZ	4:A:392:HOH:O	2.42	0.67
1:A:214:ARG:HH11	1:A:249:ILE:HG23	1.62	0.64
1:A:33:PRO:O	1:A:34:LEU:HD22	2.00	0.61
1:A:84:THR:HG22	1:A:85:ARG:H	1.66	0.59
1:A:179:ARG:HG2	1:A:179:ARG:HH11	1.74	0.52
1:A:179:ARG:HG2	1:A:179:ARG:NH1	2.24	0.52
1:A:105:ARG:HH11	1:A:105:ARG:HB3	1.74	0.52
1:A:246:GLU:H	1:A:246:GLU:CD	2.18	0.47

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A + 1	A 4 0	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}(\mathring{\rm A})$	overlap (Å)
3:A:1:UNE:O11	2:A:327:IMD:H5	2.14	0.47
1:A:269:TRP:CD1	1:A:279:PRO:HD3	2.49	0.46
1:A:212:TRP:CE3	1:A:216:HIS:HA	2.51	0.46
1:A:255:PHE:CZ	1:A:264:GLN:HG2	2.51	0.46
1:A:143:LEU:HD11	1:A:147:PHE:CZ	2.50	0.46
1:A:35:GLU:OE1	1:A:35:GLU:N	2.49	0.45
1:A:214:ARG:NE	4:A:392:HOH:O	2.49	0.45
1:A:46:SER:HA	1:A:50:GLY:O	2.17	0.45
1:A:304:LEU:O	1:A:305:HIS:CB	2.64	0.45
1:A:304:LEU:O	1:A:305:HIS:HB3	2.17	0.44
1:A:166:ARG:HH22	1:A:200:ASP:CG	2.21	0.44
1:A:214:ARG:HG2	4:A:392:HOH:O	2.17	0.43
1:A:57:ARG:HD2	1:A:60:ASP:OD1	2.18	0.43
1:A:274:ARG:HB2	4:A:20:HOH:O	2.19	0.43
1:A:174:LEU:O	1:A:182:LEU:HA	2.19	0.42
1:A:230:ILE:HG23	1:A:241:PRO:HD2	2.01	0.42
1:A:82:ASN:HD22	1:A:82:ASN:C	2.24	0.41
1:A:135:GLU:HB3	1:A:136:ARG:HG2	2.01	0.41
1:A:101:SER:HB2	1:A:183:LYS:NZ	2.35	0.41
1:A:112:ARG:HB3	1:A:113:PRO:CD	2.51	0.41
1:A:84:THR:HG22	1:A:85:ARG:N	2.34	0.41
1:A:145:ARG:HG2	1:A:293:VAL:HG23	2.03	0.41
1:A:214:ARG:NH1	1:A:249:ILE:HG23	2.32	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	272/299 (91%)	262 (96%)	7 (3%)	3 (1%)		14	30	

#### All (3) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	60	ASP
1	A	34	LEU
1	A	81	PRO

#### 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	Analysed Rotameric C		Percentiles	
1	A	245/266 (92%)	230 (94%)	15 (6%)	18 38	

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	A	76	ASP
1	A	82	ASN
1	A	84	THR
1	A	85	ARG
1	A	105	ARG
1	A	124	GLU
1	A	136	ARG
1	A	179	ARG
1	A	189	SER
1	A	239	ASP
1	A	246	GLU
1	A	258	ARG
1	A	290	MET
1	A	306	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	ASN
1	A	265	HIS
1	A	287	HIS



#### 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)

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Tuno	Chain	Res Link		Bo	nd leng	$ ag{ths}$	В	ond ang	les
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	UNE	A	1	-	11,16,16	1.46	3 (27%)	16,23,23	3.33	6 (37%)
2	IMD	A	1001	-	3,5,5	0.37	0	4,5,5	0.74	0
2	IMD	A	327	-	3,5,5	0.62	0	4,5,5	0.62	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IMD	A	327	-	-	-	0/1/1/1
2	IMD	A	1001	-	-	-	0/1/1/1
3	UNE	A	1	-	-	0/2/6/6	0/2/2/2

All (3) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
3	A	1	UNE	C09-C04	2.76	1.48	1.42
3	A	1	UNE	C04-C05	-2.66	1.37	1.43
3	A	1	UNE	O15-C13	2.10	1.36	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	A	1	UNE	O11-C06-C05	7.90	123.14	115.03
3	A	1	UNE	BR10-C08-C07	-7.38	109.02	119.27
3	A	1	UNE	C08-C09-C04	-4.71	112.70	119.32
3	A	1	UNE	C09-C08-C07	3.83	127.31	122.62
3	A	1	UNE	BR10-C08-C09	2.68	123.66	119.72
3	A	1	UNE	O14-C13-C02	-2.12	116.95	121.24

There are no chirality outliers.

There are no torsion outliers.

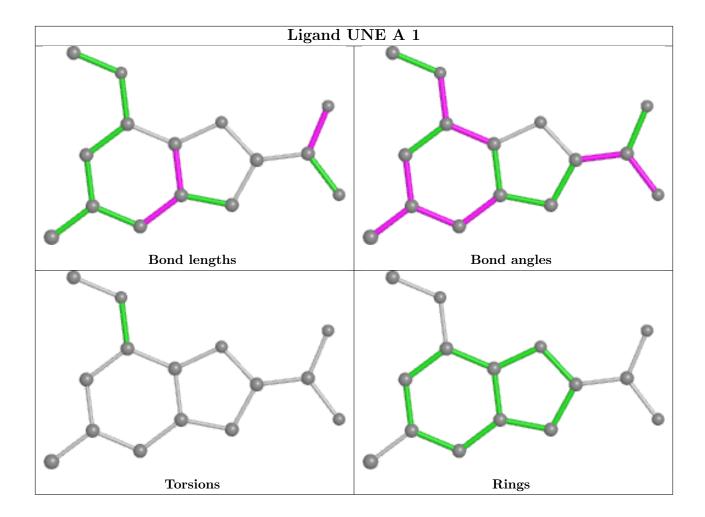
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1	UNE	1	0
2	A	327	IMD	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 then 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,  $95^{th}$  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 $>$	$\# \mathrm{RSRZ}{>}2$		$OWAB(A^2)$	Q < 0.9	
1	A	274/299 (91%)	-0.13	11 (4%)	38	31	12, 30, 63, 82	3 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	$\operatorname{Res}$	Type	RSRZ
1	A	81	PRO	6.3
1	A	80	LEU	4.8
1	A	254	PHE	4.4
1	A	82	ASN	4.3
1	A	84	THR	4.0
1	A	159	HIS	3.9
1	A	305	HIS	3.7
1	A	306	SER	3.0
1	A	83	GLY	2.4
1	A	36	SER	2.2
1	A	33	PRO	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 monosaccharides in this entry.

## 6.4 Ligands (i)

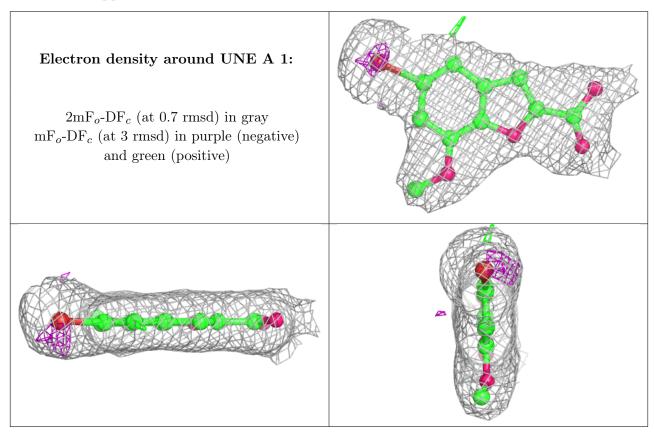
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,



median,  $95^{th}$  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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	IMD	A	327	5/5	0.94	0.18	33,35,36,37	0
2	IMD	A	1001	5/5	0.96	0.14	44,44,46,47	0
3	UNE	A	1	15/15	0.96	0.14	21,28,35,50	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.



## 6.5 Other polymers (i)

There are no such residues in this entry.

