



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 12:43 PM BST

PDB ID : 1DY9  
Title : Inhibition of the Hepatitis C Virus NS3/4A Protease. The Crystal Structures of Two Protease-Inhibitor Complexes (inhibitor I)  
Authors : Di Marco, S.; Rizzi, M.; Volpari, C.; Walsh, M.; Narjes, F.; Colarusso, S.; De Francesco, R.; Matassa, V.G.; Sollazzo, M.  
Deposited on : 2000-01-31  
Resolution : 2.10 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 : 2.13.1  
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) : 2.13.1

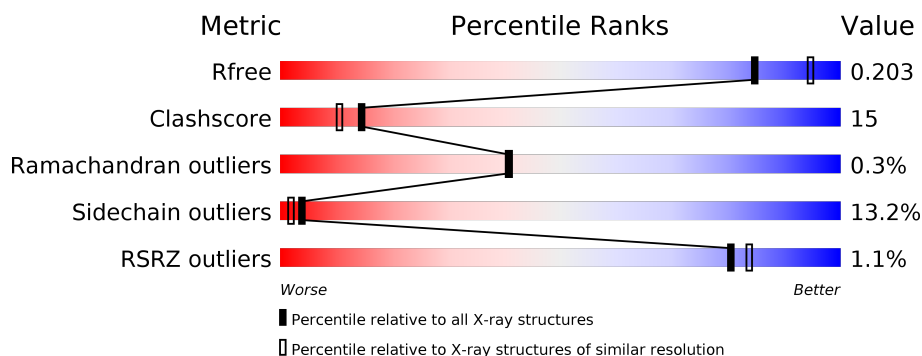
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	187	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span>%</span> <span>63%</span> <span>25%</span> <span>5%</span> <span>6%</span> </div> </div>
1	B	187	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span>%</span> <span>66%</span> <span>21%</span> <span>6%</span> <span>6%</span> </div> </div>
2	C	16	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, orange, red, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span></span> <span>56%</span> <span>6%</span> <span>6%</span> <span>6%</span> <span>25%</span> </div> </div>
2	D	16	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green, yellow, orange, red, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> <span></span> <span>50%</span> <span>19%</span> <span>6%</span> <span>25%</span> </div> </div>

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
3	2ZF	A	401	X	-	-	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3018 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 PROTEASE/HELICASE NS3 (P70).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	175	Total	C	N	O	S	0	0	0
			1283	801	231	241	10			
1	B	175	Total	C	N	O	S	0	0	0
			1283	801	231	241	10			

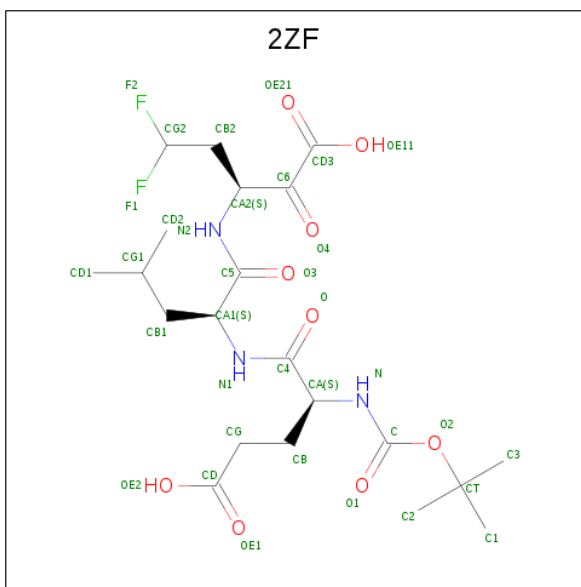
- Molecule 2 is a protein called NONSTRUCTURAL PROTEIN NS4A (P4).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	0	0	0
			84	55	15	14			
2	D	12	Total	C	N	O	0	0	0
			84	55	15	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	220	LYS	THR	engineered mutation	UNP Q81755
C	235	LYS	PRO	engineered mutation	UNP Q81755
D	220	LYS	THR	engineered mutation	UNP Q81755
D	235	LYS	PRO	engineered mutation	UNP Q81755

- Molecule 3 is N-(tert-butoxycarbonyl)-L-alpha-glutamyl-N-[(1R)-1-(carboxycarbonyl)-3,3-difluoropropyl]-L-leucinamide (three-letter code: 2ZF) (formula: C<sub>21</sub>H<sub>33</sub>F<sub>2</sub>N<sub>3</sub>O<sub>9</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			35	21	2	3	9		
3	B	1	Total	C	F	N	O	0	0
			35	21	2	3	9		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		

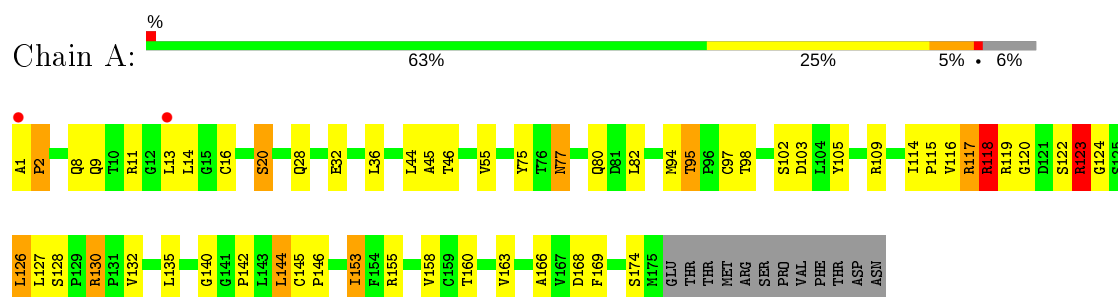
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	99	Total	O	0	0
			99	99		
5	B	103	Total	O	0	0
			103	103		
5	C	7	Total	O	0	0
			7	7		
5	D	4	Total	O	0	0
			4	4		

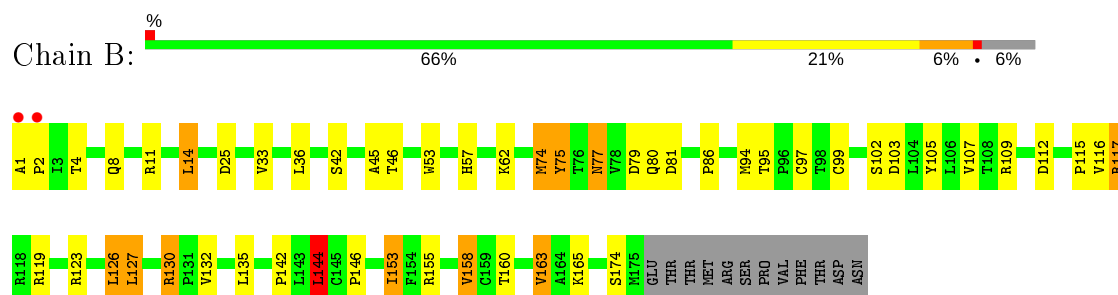
### 3 Residue-property plots [i](#)

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: PROTEASE/HELICASE NS3 (P70)



- Molecule 1: PROTEASE/HELICASE NS3 (P70)



- Molecule 2: NONSTRUCTURAL PROTEIN NS4A (P4)



- Molecule 2: NONSTRUCTURAL PROTEIN NS4A (P4)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.35Å 94.35Å 82.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.10 29.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.10) 98.2 (29.00-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.81 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.213 , 0.275 0.188 , 0.203	Depositor DCC
$R_{free}$ test set	1224 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.4	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 58.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.088 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3018	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.92% 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: 2ZF, ZN

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	0.53	0/1310	1.45	14/1786 (0.8%)
1	B	0.55	0/1310	1.33	11/1786 (0.6%)
2	C	0.57	0/83	1.42	1/111 (0.9%)
2	D	0.68	0/83	1.25	1/111 (0.9%)
All	All	0.55	0/2786	1.38	27/3794 (0.7%)

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	B	0	2

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	ARG	NE-CZ-NH2	16.53	128.56	120.30
1	A	130	ARG	NE-CZ-NH1	-16.07	112.27	120.30
1	A	118	ARG	NE-CZ-NH1	14.12	127.36	120.30
1	B	130	ARG	NE-CZ-NH2	-12.29	114.16	120.30
1	B	109	ARG	NE-CZ-NH2	10.91	125.75	120.30
1	A	117	ARG	NE-CZ-NH1	-10.46	115.07	120.30
1	B	155	ARG	NE-CZ-NH2	-9.39	115.61	120.30
1	A	123	ARG	NE-CZ-NH1	9.04	124.82	120.30
2	C	228	ARG	NE-CZ-NH2	-8.09	116.25	120.30
1	A	155	ARG	NE-CZ-NH2	-7.90	116.35	120.30
1	A	118	ARG	CD-NE-CZ	7.81	134.53	123.60
1	B	109	ARG	NE-CZ-NH1	-7.71	116.45	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	155	ARG	CD-NE-CZ	7.63	134.28	123.60
1	A	118	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	A	11	ARG	NE-CZ-NH1	7.27	123.94	120.30
1	A	109	ARG	NE-CZ-NH2	6.52	123.56	120.30
1	B	75	TYR	CB-CG-CD2	-6.49	117.11	121.00
1	A	130	ARG	CD-NE-CZ	-6.32	114.76	123.60
1	A	130	ARG	CG-CD-NE	5.92	124.24	111.80
1	B	112	ASP	CB-CG-OD2	-5.69	113.18	118.30
1	B	144	LEU	CA-CB-CG	5.66	128.31	115.30
1	B	117	ARG	NE-CZ-NH2	-5.46	117.57	120.30
2	D	228	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	55	VAL	CA-CB-CG1	-5.14	103.18	110.90
1	B	107	VAL	CA-CB-CG2	-5.05	103.33	110.90
1	B	79	ASP	CB-CG-OD2	-5.02	113.78	118.30
1	B	155	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	158	VAL	Mainchain
1	B	53	TRP	Mainchain

## 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	1283	0	1296	52	0
1	B	1283	0	1294	41	0
2	C	84	0	99	4	0
2	D	84	0	99	3	0
3	A	35	0	31	2	0
3	B	35	0	30	1	0
4	B	1	0	0	0	0
5	A	99	0	0	3	0
5	B	103	0	0	3	0
5	C	7	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	4	0	0	0	0
All	All	3018	0	2849	87	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 15.

All (87) 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:A:117:ARG:NH1	1:B:115:PRO:HB3	1.76	1.01
1:A:123:ARG:HG3	1:A:123:ARG:HH11	1.39	0.87
1:B:74:MET:HE3	1:B:75:TYR:HB2	1.58	0.83
1:A:8:GLN:HE21	2:C:228:ARG:HE	1.28	0.82
1:B:74:MET:CE	1:B:75:TYR:HB2	2.11	0.81
1:B:80:GLN:OE1	1:B:174:SER:HB3	1.81	0.80
1:B:75:TYR:OH	1:B:80:GLN:NE2	2.15	0.79
1:A:80:GLN:OE1	1:A:174:SER:HB3	1.86	0.76
1:B:77:ASN:CG	1:B:80:GLN:HG2	2.06	0.75
1:A:75:TYR:OH	1:A:80:GLN:NE2	2.20	0.74
1:A:117:ARG:HH12	1:B:115:PRO:HB3	1.51	0.72
1:B:46:THR:HG23	1:B:94:MET:HE1	1.73	0.70
3:A:401:2ZF:CB2	3:A:401:2ZF:OE11	2.35	0.69
1:B:4:THR:OG1	2:D:232:SER:HB2	1.94	0.68
1:B:46:THR:HG23	1:B:94:MET:CE	2.24	0.68
1:B:163:VAL:HG22	1:B:165:LYS:HE3	1.78	0.66
1:B:8:GLN:HE21	2:D:228:ARG:HE	1.43	0.65
1:A:142:PRO:HB2	1:A:144:LEU:HD13	1.79	0.63
1:A:46:THR:HG23	1:A:94:MET:CE	2.29	0.62
1:A:123:ARG:CG	1:A:123:ARG:HH11	2.11	0.62
1:B:142:PRO:HB2	1:B:144:LEU:HD13	1.81	0.62
1:A:77:ASN:CB	1:A:80:GLN:HE21	2.13	0.61
1:A:77:ASN:C	1:A:77:ASN:HD22	2.05	0.61
1:A:119:ARG:O	1:B:117:ARG:NH2	2.34	0.60
3:A:401:2ZF:HB1	3:A:401:2ZF:OE11	1.71	0.60
1:A:117:ARG:HH12	1:B:115:PRO:CB	2.13	0.60
1:B:46:THR:HB	1:B:153:ILE:HD11	1.85	0.58
1:B:57:HIS:ND1	1:B:81:ASP:OD2	2.33	0.57
1:A:80:GLN:HG3	1:A:82:LEU:HB3	1.87	0.57
1:B:11:ARG:NH2	1:B:25:ASP:OD1	2.38	0.57
1:A:117:ARG:NH1	1:B:115:PRO:CB	2.60	0.56
1:A:97:CYS:O	1:A:98:THR:HB	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:VAL:HG22	1:A:126:LEU:HD12	1.87	0.56
3:B:401:2ZF:HB22	3:B:401:2ZF:OE11	2.05	0.55
1:A:46:THR:HG23	1:A:94:MET:HE1	1.88	0.55
1:A:119:ARG:HH22	1:B:102:SER:HB2	1.72	0.55
1:A:8:GLN:NE2	2:C:228:ARG:HE	2.01	0.55
1:A:142:PRO:HB2	1:A:144:LEU:CD1	2.37	0.55
1:A:16:CYS:SG	5:A:2017:HOH:O	2.58	0.54
1:B:74:MET:HE2	1:B:75:TYR:HB2	1.91	0.52
1:A:16:CYS:O	1:A:20:SER:OG	2.29	0.51
1:B:163:VAL:CG2	1:B:165:LYS:HE3	2.40	0.51
1:A:9:GLN:HG3	5:C:2004:HOH:O	2.10	0.51
1:A:117:ARG:HD2	1:B:127:LEU:HD11	1.93	0.50
1:B:2:PRO:HD2	1:B:105:TYR:CZ	2.46	0.50
1:A:77:ASN:CG	1:A:80:GLN:HG2	2.33	0.49
1:A:95:THR:HG22	5:A:2034:HOH:O	2.12	0.49
1:A:8:GLN:HE22	2:C:228:ARG:HH21	1.61	0.49
1:B:103:ASP:O	1:B:146:PRO:HD3	2.14	0.48
1:B:11:ARG:HH21	1:B:25:ASP:CG	2.16	0.48
1:B:62:LYS:HE2	5:B:2037:HOH:O	2.13	0.48
1:A:118:ARG:HD2	1:A:120:GLY:O	2.13	0.48
1:A:77:ASN:ND2	1:A:80:GLN:H	2.12	0.47
1:A:77:ASN:HB3	1:A:80:GLN:HG2	1.95	0.47
1:A:119:ARG:HH22	1:B:102:SER:CB	2.27	0.47
1:A:46:THR:HB	1:A:153:ILE:HD11	1.96	0.47
1:B:1:ALA:HB3	5:B:2001:HOH:O	2.14	0.47
1:A:77:ASN:CG	1:A:80:GLN:HE21	2.18	0.46
1:B:33:VAL:HB	2:D:229:ILE:HB	1.96	0.46
1:A:123:ARG:CG	1:A:123:ARG:NH1	2.74	0.45
1:B:36:LEU:HD11	1:B:45:ALA:HB2	1.97	0.45
1:A:2:PRO:HD2	1:A:105:TYR:CE2	2.52	0.45
1:A:36:LEU:HD11	1:A:45:ALA:HB2	1.99	0.45
1:A:124:GLY:O	1:A:166:ALA:HB1	2.18	0.44
1:B:132:VAL:HG12	5:B:2091:HOH:O	2.18	0.43
1:B:46:THR:CB	1:B:153:ILE:HD11	2.48	0.43
1:A:13:LEU:HD23	1:A:13:LEU:HA	1.89	0.43
1:A:77:ASN:OD1	1:A:80:GLN:NE2	2.52	0.43
1:B:142:PRO:HB2	1:B:144:LEU:CD1	2.48	0.43
1:B:46:THR:CG2	1:B:94:MET:HE1	2.46	0.43
1:B:116:VAL:HG22	1:B:126:LEU:CD1	2.49	0.42
1:A:130:ARG:NH2	5:A:2079:HOH:O	2.51	0.42
1:A:32:GLU:OE1	2:C:231:LEU:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ARG:HG3	1:A:123:ARG:NH1	2.16	0.42
1:B:77:ASN:C	1:B:77:ASN:HD22	2.23	0.42
1:A:122:SER:HB2	1:A:169:PHE:O	2.20	0.42
1:B:14:LEU:HA	1:B:14:LEU:HD23	1.94	0.41
1:B:130:ARG:HA	1:B:130:ARG:HD2	1.91	0.41
1:A:145:CYS:HB2	1:A:146:PRO:HD2	2.02	0.41
1:A:102:SER:CB	1:B:127:LEU:HD12	2.51	0.41
1:A:127:LEU:HD23	1:A:127:LEU:HA	1.84	0.41
1:A:1:ALA:HA	1:A:2:PRO:HD2	1.97	0.41
1:A:114:ILE:HA	1:A:115:PRO:HD3	1.79	0.41
1:B:74:MET:HB3	1:B:86:PRO:HD3	2.03	0.41
1:A:77:ASN:CB	1:A:80:GLN:HG2	2.51	0.40
1:A:80:GLN:HG3	1:A:82:LEU:CB	2.51	0.40
1:A:44:LEU:O	1:A:140:GLY:HA3	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	173/187 (92%)	168 (97%)	4 (2%)	1 (1%)	25	21
1	B	173/187 (92%)	173 (100%)	0	0	100	100
2	C	10/16 (62%)	10 (100%)	0	0	100	100
2	D	10/16 (62%)	10 (100%)	0	0	100	100
All	All	366/406 (90%)	361 (99%)	4 (1%)	1 (0%)	41	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	PRO

### 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	142/154 (92%)	124 (87%)	18 (13%)	4	2
1	B	142/154 (92%)	125 (88%)	17 (12%)	5	2
2	C	10/13 (77%)	7 (70%)	3 (30%)	0	0
2	D	10/13 (77%)	8 (80%)	2 (20%)	1	0
All	All	304/334 (91%)	264 (87%)	40 (13%)	4	2

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	20	SER
1	A	28	GLN
1	A	77	ASN
1	A	95	THR
1	A	103	ASP
1	A	118	ARG
1	A	123	ARG
1	A	126	LEU
1	A	128	SER
1	A	132	VAL
1	A	135	LEU
1	A	144	LEU
1	A	153	ILE
1	A	158	VAL
1	A	160	THR
1	A	163	VAL
1	A	168	ASP
1	B	14	LEU
1	B	42	SER
1	B	74	MET
1	B	77	ASN
1	B	95	THR
1	B	97	CYS

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Mol	Chain	Res	Type
1	B	99	CYS
1	B	119	ARG
1	B	123	ARG
1	B	126	LEU
1	B	127	LEU
1	B	135	LEU
1	B	144	LEU
1	B	153	ILE
1	B	158	VAL
1	B	160	THR
1	B	163	VAL
2	C	228	ARG
2	C	231	LEU
2	C	232	SER
2	D	228	ARG
2	D	231	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	27	ASN
1	A	28	GLN
1	A	34	GLN
1	A	77	ASN
1	A	80	GLN
1	B	8	GLN
1	B	27	ASN
1	B	34	GLN
1	B	77	ASN
1	B	80	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 3 ligands modelled in this entry, 1 is 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	2ZF	A	401	1	28,34,34	3.57	5 (17%)	34,47,47	1.63	4 (11%)
3	2ZF	B	401	1	28,34,34	3.45	5 (17%)	34,47,47	2.34	6 (17%)

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	2ZF	A	401	1	1/1/12/15	8/40/46/46	-
3	2ZF	B	401	1	-	10/40/46/46	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	2ZF	O4-C6	17.28	1.46	1.22
3	B	401	2ZF	O4-C6	16.83	1.45	1.22
3	A	401	2ZF	CA2-C6	4.79	1.61	1.53
3	B	401	2ZF	CA2-C6	3.64	1.59	1.53
3	B	401	2ZF	CB2-CG2	3.40	1.56	1.50
3	A	401	2ZF	CB2-CG2	3.37	1.56	1.50
3	A	401	2ZF	O2-CT	-2.71	1.43	1.48
3	B	401	2ZF	C4-N1	2.34	1.39	1.34
3	B	401	2ZF	O2-CT	-2.26	1.44	1.48
3	A	401	2ZF	C4-N1	2.07	1.38	1.34

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	2ZF	CB2-CA2-N2	8.99	131.28	110.58
3	A	401	2ZF	CB2-CA2-N2	6.42	125.37	110.58
3	B	401	2ZF	CB-CA-N	5.19	121.37	110.88
3	B	401	2ZF	C6-CA2-N2	4.69	116.56	109.85
3	B	401	2ZF	CG-CB-CA	-3.92	105.49	113.21
3	A	401	2ZF	CA2-N2-C5	3.13	128.37	121.67
3	B	401	2ZF	C5-CA1-N1	-3.07	102.81	111.16
3	A	401	2ZF	C5-CA1-N1	-2.75	103.66	111.16
3	A	401	2ZF	CB-CA-N	2.69	116.31	110.88
3	B	401	2ZF	CB-CG-CD	-2.14	108.99	113.59

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	401	2ZF	CA2

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	401	2ZF	N-CA-CB-CG
3	A	401	2ZF	C4-CA-CB-CG
3	A	401	2ZF	CB2-CA2-N2-C5
3	A	401	2ZF	CD3-C6-CA2-CB2
3	A	401	2ZF	CA2-CB2-CG2-F1
3	A	401	2ZF	CA2-CB2-CG2-F2
3	B	401	2ZF	O4-C6-CA2-CB2
3	B	401	2ZF	CD3-C6-CA2-CB2
3	B	401	2ZF	N2-CA2-CB2-CG2
3	B	401	2ZF	CA2-CB2-CG2-F1
3	B	401	2ZF	CA-CB-CG-CD
3	B	401	2ZF	O-C4-N1-CA1
3	A	401	2ZF	C6-CA2-CB2-CG2
3	B	401	2ZF	CA2-CB2-CG2-F2
3	A	401	2ZF	N2-CA2-CB2-CG2
3	B	401	2ZF	CA1-CB1-CG1-CD2
3	B	401	2ZF	O3-C5-N2-CA2
3	B	401	2ZF	O-C4-CA-N

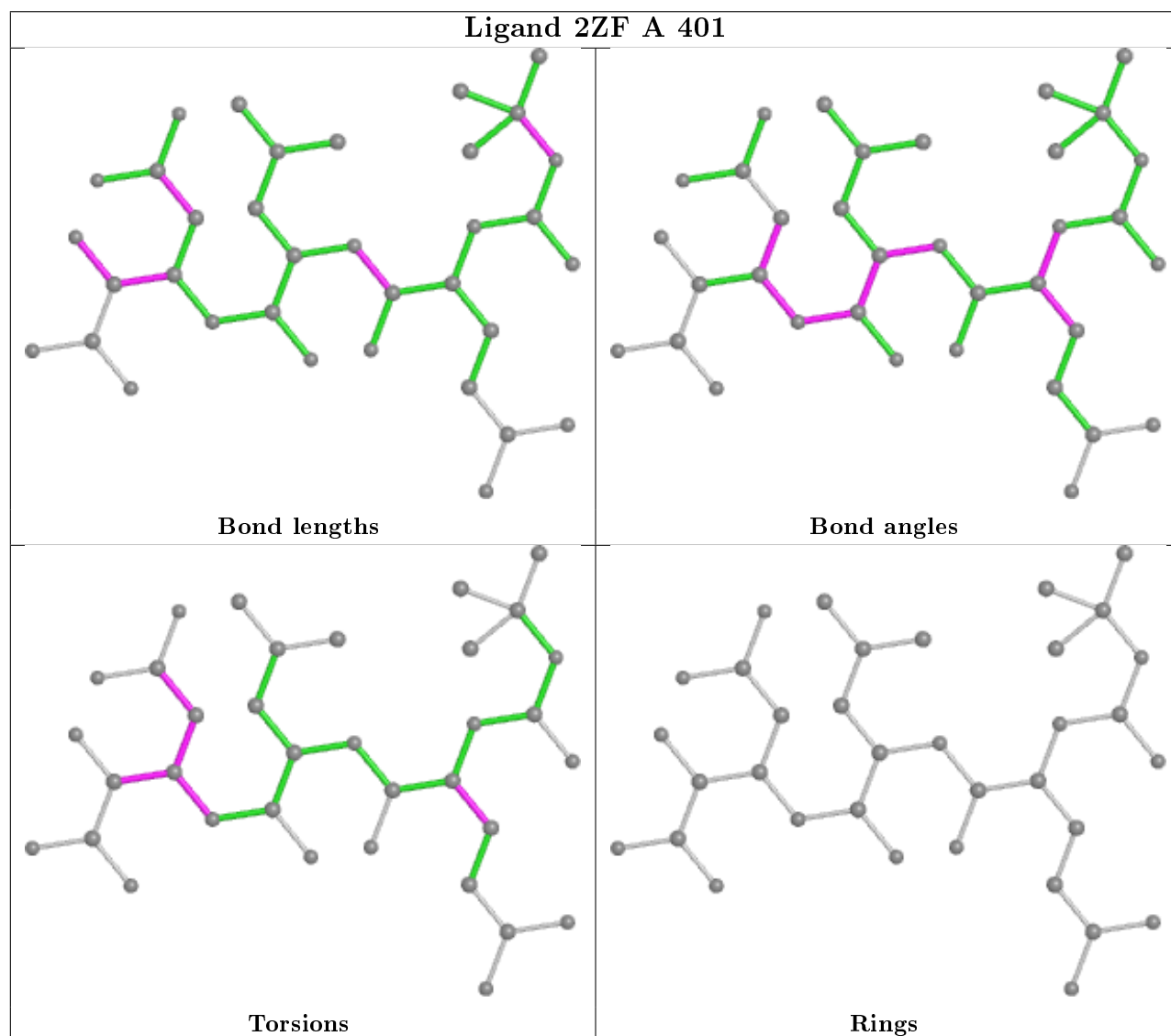
There are no ring outliers.

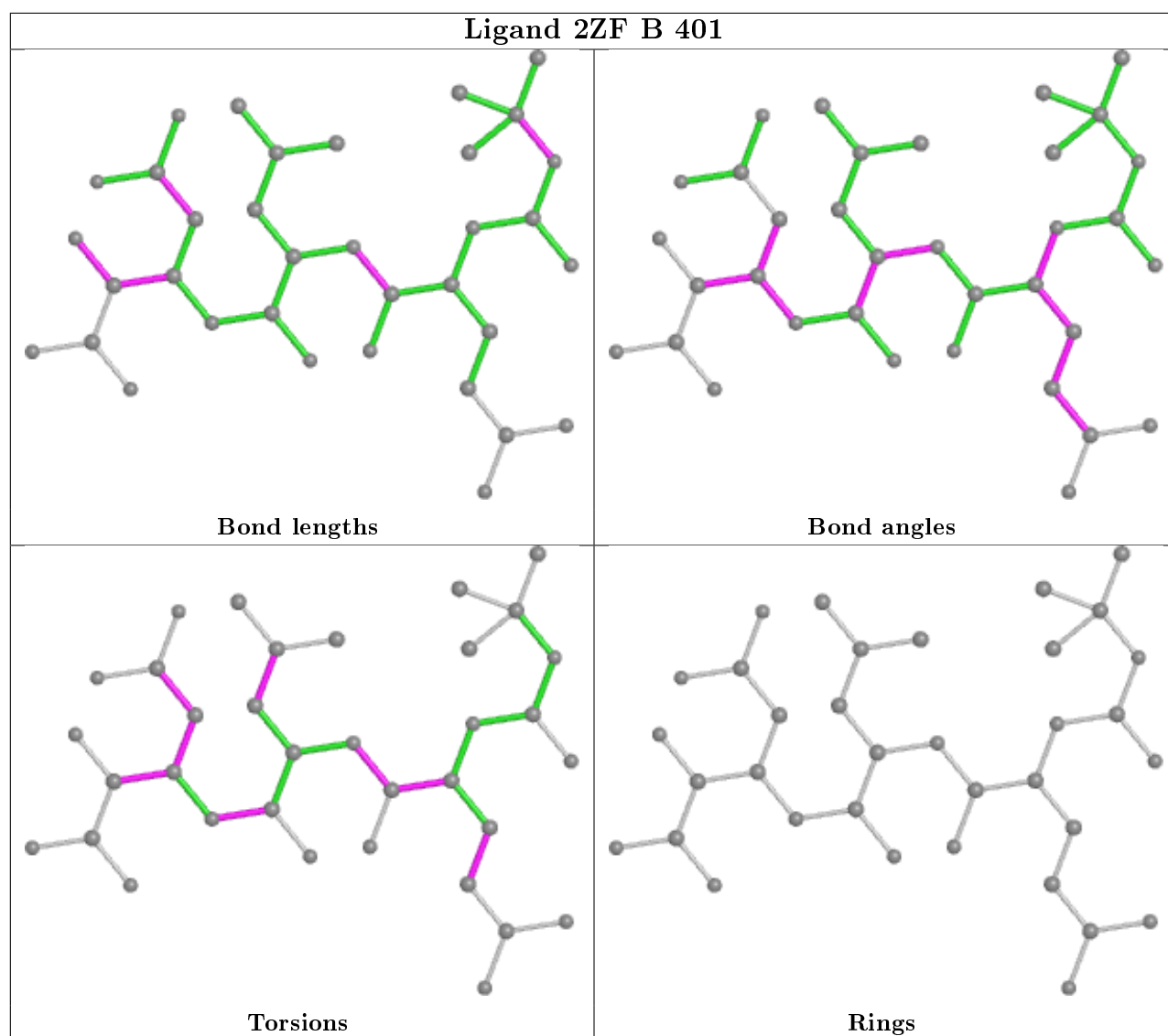
2 monomers are involved in 3 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	2ZF	2	0
3	B	401	2ZF	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	175/187 (93%)	-0.14	2 (1%) 80 84	25, 41, 68, 93	0
1	B	175/187 (93%)	-0.18	2 (1%) 80 84	25, 39, 61, 86	0
2	C	12/16 (75%)	-0.10	0 100 100	28, 33, 47, 61	0
2	D	12/16 (75%)	-0.49	0 100 100	27, 31, 40, 57	0
All	All	374/406 (92%)	-0.17	4 (1%) 80 84	25, 40, 66, 93	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	ALA	7.6
1	A	1	ALA	3.8
1	A	13	LEU	3.0
1	B	2	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](#)

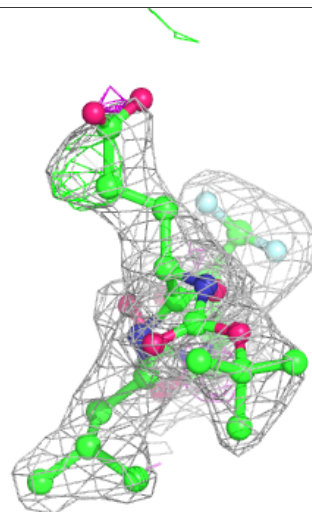
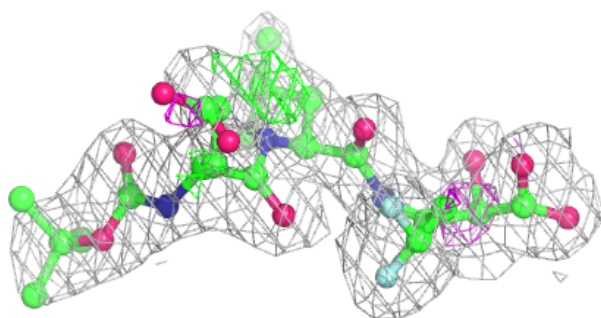
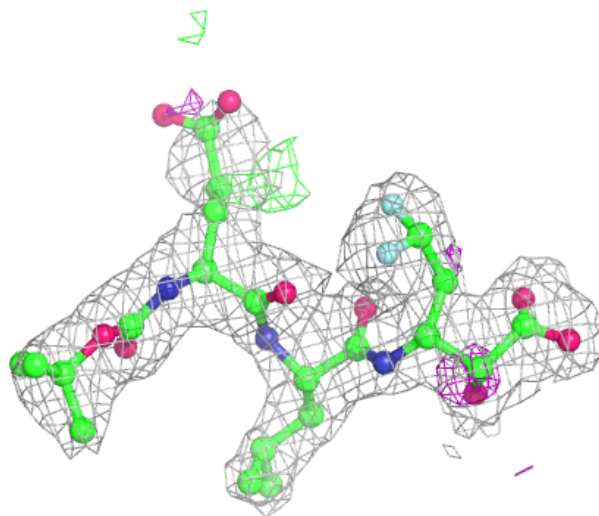
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	2ZF	A	401	35/35	0.89	0.17	41,50,76,80	0
4	ZN	B	300	1/1	0.93	0.07	81,81,81,81	0
3	2ZF	B	401	35/35	0.94	0.11	38,46,60,61	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.

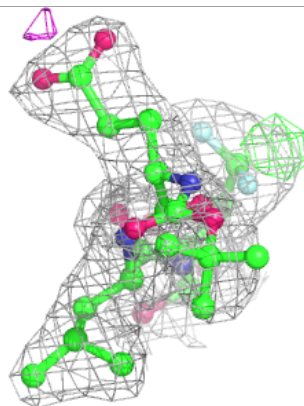
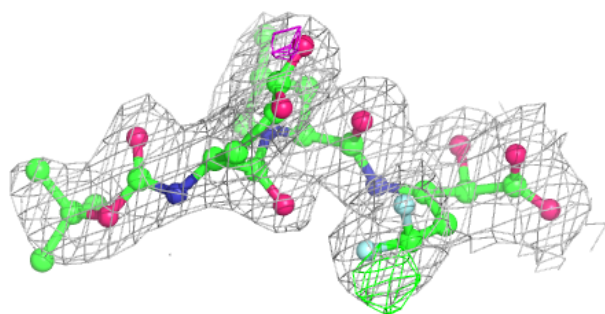
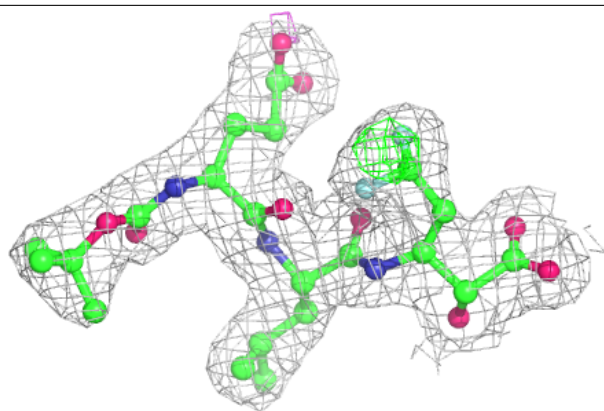
**Electron density around 2ZF A 401:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around 2ZF B 401:**

$2mF_o-DF_c$  (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.