



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2021 – 12:45 PM EDT

PDB ID : 7JXL  
Title : EGFR kinase (T790M/V948R) in complex with AZ5104  
Authors : Beyett, T.S.; Eck, M.J.  
Deposited on : 2020-08-27  
Resolution : 2.40 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 : 2.23.1  
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) : 2.23.1

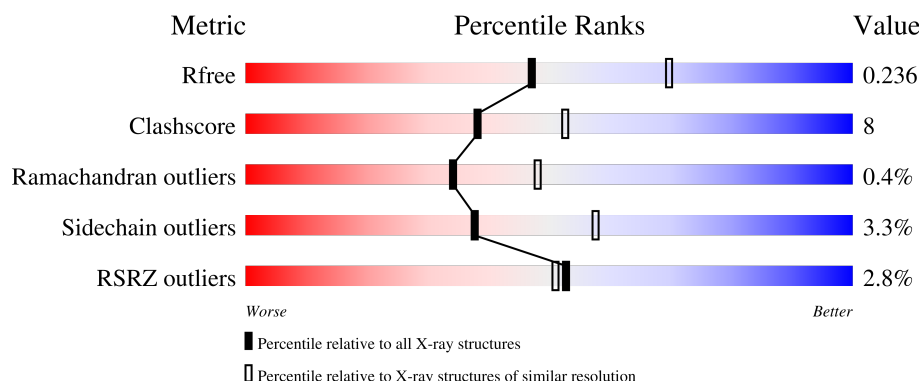
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	331	 2% 76% 12% • 11%
1	B	331	 3% 72% 13% • 14%
1	C	331	 3% 72% 15% • 12%
1	D	331	 2% 69% 18% • 12%

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9668 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.

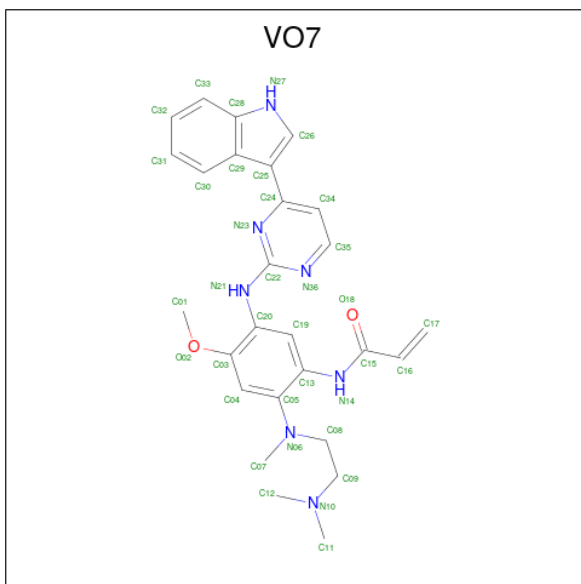
- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	292	Total	C	N	O	S	0	1	0
			2362	1517	402	423	20			
1	A	294	Total	C	N	O	S	0	1	0
			2378	1528	402	428	20			
1	C	292	Total	C	N	O	S	0	1	0
			2365	1521	402	422	20			
1	B	285	Total	C	N	O	S	0	1	0
			2300	1480	390	412	18			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	692	GLY	-	expression tag	UNP P00533
D	693	SER	-	expression tag	UNP P00533
D	694	THR	-	expression tag	UNP P00533
D	790	MET	THR	engineered mutation	UNP P00533
D	948	ARG	VAL	engineered mutation	UNP P00533
A	692	GLY	-	expression tag	UNP P00533
A	693	SER	-	expression tag	UNP P00533
A	694	THR	-	expression tag	UNP P00533
A	790	MET	THR	engineered mutation	UNP P00533
A	948	ARG	VAL	engineered mutation	UNP P00533
C	692	GLY	-	expression tag	UNP P00533
C	693	SER	-	expression tag	UNP P00533
C	694	THR	-	expression tag	UNP P00533
C	790	MET	THR	engineered mutation	UNP P00533
C	948	ARG	VAL	engineered mutation	UNP P00533
B	692	GLY	-	expression tag	UNP P00533
B	693	SER	-	expression tag	UNP P00533
B	694	THR	-	expression tag	UNP P00533
B	790	MET	THR	engineered mutation	UNP P00533
B	948	ARG	VAL	engineered mutation	UNP P00533

- Molecule 2 is N-(2-([2-(dimethylamino)ethyl](methyl)amino}-5-([4-(1H-indol-3-yl)pyrimidin-2-yl]amino}-4-methoxyphenyl)propanamide (three-letter code: VO7) (formula: C<sub>27</sub>H<sub>31</sub>N<sub>7</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			36	27	7	2		
2	A	1	Total	C	N	O	0	0
			36	27	7	2		
2	C	1	Total	C	N	O	0	0
			36	27	7	2		
2	B	1	Total	C	N	O	0	0
			36	27	7	2		

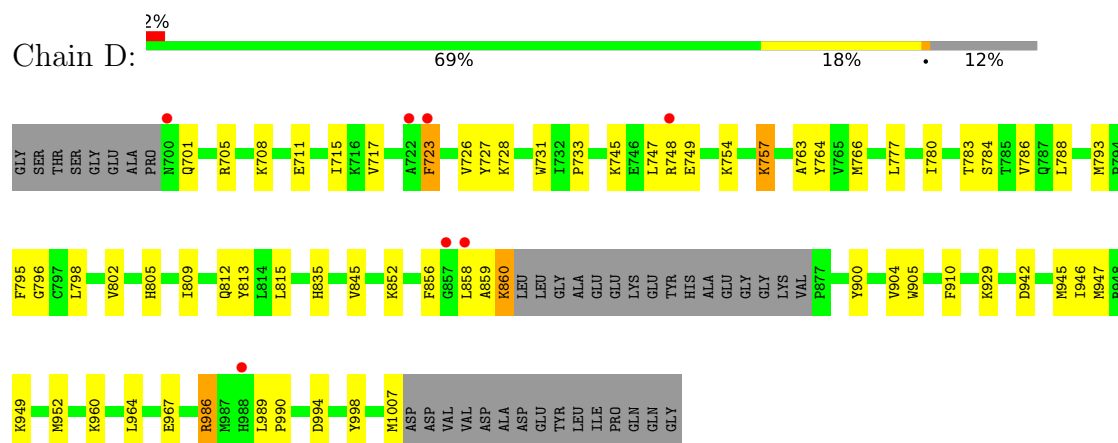
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	37	Total	O	0	0
			37	37		
3	A	37	Total	O	0	0
			37	37		
3	C	19	Total	O	0	0
			19	19		
3	B	26	Total	O	0	0
			26	26		

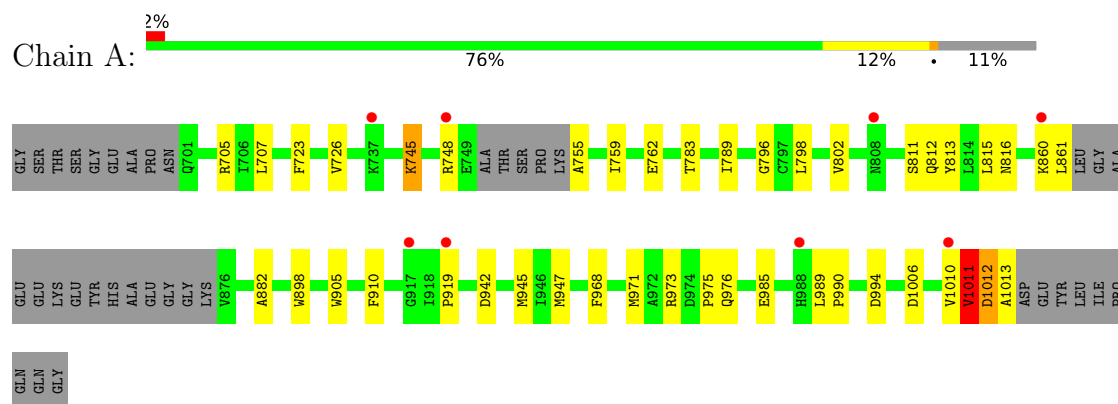
### 3 Residue-property plots [i](#)

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.

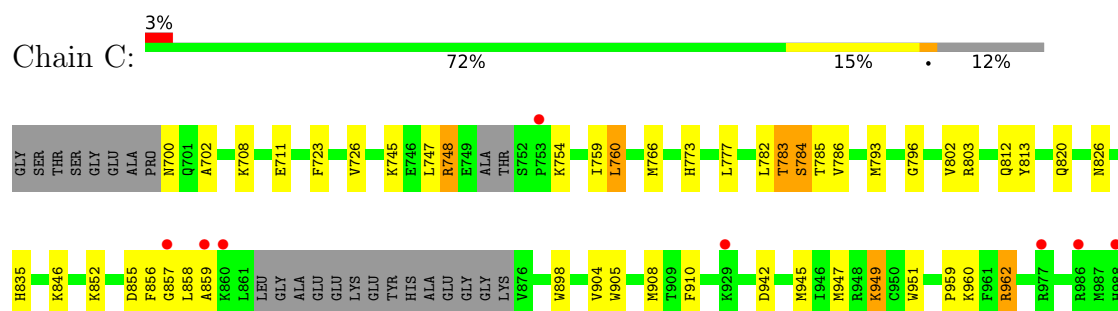
- Molecule 1: Epidermal growth factor receptor

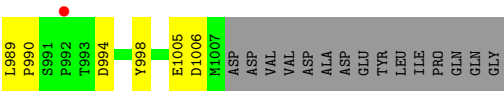


- Molecule 1: Epidermal growth factor receptor

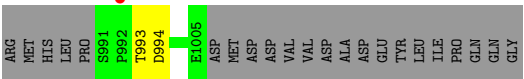
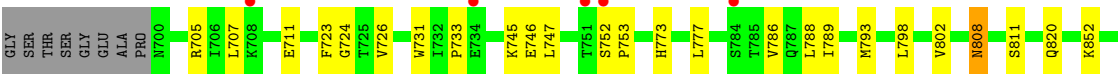


- Molecule 1: Epidermal growth factor receptor





● Molecule 1: Epidermal growth factor receptor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.05Å 101.33Å 86.95Å 90.00° 101.36° 90.00°	Depositor
Resolution (Å)	60.05 – 2.40 60.05 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (60.05-2.40) 99.6 (60.05-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.40Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.209 , 0.239 0.208 , 0.236	Depositor DCC
$R_{free}$ test set	2307 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.1	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 38.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9668	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.67% 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: VO7

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	0.46	0/2431	0.64	0/3285
1	B	0.48	0/2352	0.68	0/3178
1	C	0.45	0/2419	0.63	0/3267
1	D	0.47	0/2417	0.67	1/3265 (0.0%)
All	All	0.47	0/9619	0.66	1/12995 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	858	LEU	CB-CG-CD2	5.33	120.06	111.00

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	2378	0	2420	30	0
1	B	2300	0	2351	38	0
1	C	2365	0	2416	41	0
1	D	2362	0	2410	49	0
2	A	36	0	0	1	0
2	B	36	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	36	0	0	3	0
2	D	36	0	0	2	0
3	A	37	0	0	0	0
3	B	26	0	0	1	0
3	C	19	0	0	3	0
3	D	37	0	0	0	0
All	All	9668	0	9597	146	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 8.

All (146) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:945:MET:CE	1:B:918:ILE:CD1	2.44	0.95
1:D:945:MET:CE	1:B:918:ILE:HD11	2.04	0.87
1:A:919:PRO:HG2	1:C:945:MET:SD	2.15	0.86
1:D:945:MET:HE1	1:B:918:ILE:CD1	2.08	0.83
1:D:945:MET:HE1	1:B:918:ILE:HD11	1.60	0.82
1:C:723:PHE:HE2	1:C:745:LYS:HB3	1.54	0.72
1:A:816:ASN:ND2	1:A:1010:VAL:HG13	2.05	0.71
1:A:812:GLN:HG2	1:A:989:LEU:HG	1.73	0.70
1:B:723:PHE:HE2	1:B:745:LYS:HB3	1.57	0.70
1:B:808:ASN:O	1:B:808:ASN:ND2	2.26	0.68
1:D:705:ARG:NH2	1:D:711:GLU:OE2	2.26	0.68
1:A:905:TRP:HD1	1:A:947:MET:HE1	1.61	0.66
1:D:905:TRP:HD1	1:D:947:MET:HE1	1.61	0.66
1:A:802:VAL:HG12	1:A:910:PHE:HA	1.77	0.65
1:C:846:LYS:NZ	1:C:1005:GLU:OE2	2.27	0.65
1:C:812:GLN:HG2	1:C:989:LEU:HG	1.80	0.64
1:C:962:ARG:HG2	3:C:1205:HOH:O	1.97	0.64
1:D:945:MET:HE2	1:B:918:ILE:HD11	1.80	0.63
1:C:783:THR:OG1	1:C:784:SER:N	2.31	0.63
1:D:748:ARG:HG3	1:D:749:GLU:HG3	1.81	0.62
1:A:976:GLN:HE22	1:A:985:GLU:HB3	1.65	0.62
1:D:945:MET:HE2	1:B:918:ILE:CD1	2.27	0.62
1:A:1012:ASP:OD1	1:A:1012:ASP:N	2.22	0.62
1:A:762:GLU:OE1	1:A:860:LYS:HB2	1.99	0.62
1:D:835:HIS:CD2	1:D:856:PHE:HB3	2.35	0.61
1:D:723:PHE:HD1	1:D:723:PHE:H	1.47	0.61
1:C:949:LYS:HB3	1:C:959:PRO:HD3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:796:GLY:HA2	2:A:1101:VO7:C19	2.33	0.58
1:A:802:VAL:CG1	1:A:910:PHE:HA	2.33	0.58
1:B:802:VAL:HG12	1:B:910:PHE:HA	1.83	0.58
1:D:802:VAL:HG12	1:D:910:PHE:HA	1.84	0.58
1:C:783:THR:HG23	1:C:785:THR:O	2.02	0.58
1:A:816:ASN:ND2	1:A:1011:VAL:HG23	2.20	0.57
1:C:962:ARG:HB3	1:C:962:ARG:HH21	1.69	0.57
1:A:811:SER:OG	1:A:975:PRO:HB2	2.06	0.56
1:C:962:ARG:HB3	1:C:962:ARG:NH2	2.20	0.56
1:C:759:ILE:HD13	1:C:786:VAL:HG21	1.88	0.56
1:A:905:TRP:CD1	1:A:947:MET:HE1	2.41	0.56
1:D:942:ASP:OD2	1:B:929:LYS:HE2	2.05	0.56
1:B:705:ARG:NH2	1:B:711:GLU:OE2	2.39	0.56
1:D:929:LYS:HE2	1:A:942:ASP:OD2	2.06	0.55
1:C:857:GLY:O	1:C:859:ALA:N	2.39	0.55
1:B:882:ALA:HA	1:B:898:TRP:CD2	2.41	0.55
1:B:798:LEU:O	1:B:802:VAL:HG23	2.06	0.55
1:C:905:TRP:HD1	1:C:947:MET:HE1	1.72	0.55
1:D:802:VAL:CG1	1:D:910:PHE:HA	2.36	0.55
1:A:723:PHE:HB3	1:A:861:LEU:HD23	1.89	0.55
1:C:760:LEU:HD11	1:C:782:LEU:HD11	1.88	0.55
1:C:835:HIS:CD2	1:C:856:PHE:HB3	2.42	0.55
1:D:812:GLN:HG2	1:D:989:LEU:HG	1.88	0.54
1:D:990:PRO:HB2	1:D:994:ASP:HB2	1.90	0.54
1:C:802:VAL:CG1	1:C:910:PHE:HA	2.38	0.54
1:B:904:VAL:O	1:B:908[B]:MET:HG3	2.08	0.54
1:C:723:PHE:CE2	1:C:745:LYS:HD2	2.44	0.53
1:D:945:MET:CE	1:B:918:ILE:HD12	2.38	0.52
1:B:802:VAL:CG1	1:B:910:PHE:HA	2.39	0.52
1:B:747:LEU:HD12	1:B:786:VAL:HB	1.91	0.52
1:C:708:LYS:O	1:C:711:GLU:HG2	2.10	0.51
1:C:802:VAL:HG12	1:C:910:PHE:HA	1.90	0.51
1:C:723:PHE:CE2	1:C:745:LYS:HB3	2.41	0.51
1:D:945:MET:HE1	1:B:918:ILE:HD12	1.91	0.51
1:C:813:TYR:OH	1:C:990:PRO:HD3	2.11	0.50
1:B:967:GLU:HG2	1:B:971:MET:HE2	1.92	0.50
1:D:708:LYS:O	1:D:711:GLU:HG2	2.12	0.49
1:D:989:LEU:HD22	1:D:990:PRO:HD2	1.95	0.49
1:A:813:TYR:OH	1:A:990:PRO:HD3	2.12	0.49
1:D:798:LEU:O	1:D:802:VAL:HG23	2.12	0.49
1:D:793:MET:CE	1:D:852:LYS:HD3	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:816:ASN:HD21	1:A:1011:VAL:HG23	1.78	0.48
1:D:763:ALA:HA	1:D:766:MET:HE3	1.94	0.48
1:C:904:VAL:O	1:C:908[B]:MET:HG3	2.13	0.48
1:A:798:LEU:O	1:A:802:VAL:HG23	2.13	0.48
1:B:752:SER:HB2	1:B:753:PRO:HD2	1.95	0.48
1:A:726:VAL:HG22	1:A:745:LYS:HG2	1.95	0.48
1:C:796:GLY:HA2	2:C:1101:VO7:C19	2.44	0.47
1:B:793:MET:CE	1:B:852:LYS:HD3	2.45	0.47
1:A:919:PRO:HG2	1:C:945:MET:CE	2.44	0.47
1:C:723:PHE:H	1:C:723:PHE:HD1	1.62	0.47
1:D:946:ILE:HD11	1:D:967:GLU:HG2	1.97	0.47
1:A:882:ALA:HA	1:A:898:TRP:CD2	2.49	0.47
1:A:973:ARG:HA	1:A:1013:ALA:HB3	1.96	0.46
1:C:949:LYS:HD2	1:C:959:PRO:HD3	1.97	0.46
1:D:726:VAL:HG21	2:D:1101:VO7:C28	2.46	0.46
1:D:815:LEU:HD23	1:D:815:LEU:HA	1.77	0.46
1:C:826:ASN:ND2	3:C:1205:HOH:O	2.48	0.46
1:D:949:LYS:HG2	1:D:952:MET:HE1	1.99	0.45
1:B:922:GLU:O	1:B:926:ILE:HG23	2.16	0.45
1:A:990:PRO:HB2	1:A:994:ASP:HB2	1.98	0.45
1:D:793:MET:HE1	1:D:852:LYS:HD3	1.98	0.45
1:A:968:PHE:HA	1:A:971:MET:HE3	1.97	0.45
1:C:726:VAL:HG21	2:C:1101:VO7:C28	2.46	0.45
1:D:757:LYS:C	1:D:757:LYS:HD3	2.37	0.45
1:D:780:ILE:HG13	1:D:788:LEU:HD23	1.98	0.45
1:B:705:ARG:HE	1:B:707:LEU:HD23	1.82	0.45
1:D:813:TYR:OH	1:D:990:PRO:HD3	2.18	0.44
1:B:968:PHE:HA	1:B:971:MET:HE3	1.99	0.44
1:D:731:TRP:CH2	1:D:733:PRO:HG3	2.52	0.44
1:A:755:ALA:O	1:A:759:ILE:HG13	2.18	0.44
1:C:905:TRP:CD1	1:C:947:MET:HE1	2.52	0.44
1:C:700:ASN:OD1	1:C:702:ALA:HB3	2.18	0.44
1:B:994:ASP:N	1:B:994:ASP:OD1	2.49	0.44
1:B:908[B]:MET:SD	1:B:939:CYS:SG	3.16	0.44
1:B:968:PHE:CD1	1:B:971:MET:HE3	2.52	0.44
1:D:813:TYR:CE2	1:D:989:LEU:HD23	2.52	0.44
1:C:773:HIS:CD2	1:C:820:GLN:HB3	2.53	0.44
1:D:717:VAL:HG22	1:D:727:TYR:CE2	2.53	0.43
1:B:777:LEU:HD11	1:B:788:LEU:HB3	1.99	0.43
1:A:707:LEU:HD12	1:A:789:ILE:HD13	2.01	0.43
1:D:795:PHE:HB2	1:D:845:VAL:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:726:VAL:HG21	2:B:1101:VO7:C28	2.48	0.43
1:D:747:LEU:HD12	1:D:786:VAL:HB	2.01	0.42
1:C:855:ASP:OD2	2:C:1101:VO7:N27	2.53	0.42
1:D:715:ILE:HG12	1:D:728:LYS:O	2.20	0.42
1:B:773:HIS:CD2	1:B:820:GLN:HB3	2.54	0.42
1:B:811:SER:OG	1:B:975:PRO:HB2	2.20	0.42
1:D:905:TRP:CD1	1:D:947:MET:HE1	2.46	0.42
1:C:898:TRP:CE3	1:C:951:TRP:HA	2.54	0.42
1:C:748:ARG:H	1:C:748:ARG:HG3	1.49	0.42
1:C:803:ARG:NH1	3:C:1207:HOH:O	2.51	0.42
1:B:707:LEU:HD12	1:B:789:ILE:HD13	2.02	0.42
1:D:701:GLN:HA	1:D:764:TYR:CE1	2.55	0.42
1:A:968:PHE:CD1	1:A:971:MET:HE3	2.55	0.42
1:B:993:THR:O	3:B:1201:HOH:O	2.22	0.42
1:D:905:TRP:HB2	1:D:947:MET:HE3	2.02	0.42
1:D:796:GLY:HA2	2:D:1101:VO7:C19	2.49	0.41
1:D:900:TYR:CE2	1:D:964:LEU:HD13	2.55	0.41
1:C:766:MET:HE3	1:C:777:LEU:HD22	2.01	0.41
1:B:724:GLY:HA3	1:B:746:GLU:O	2.20	0.41
1:A:815:LEU:HD12	1:A:975:PRO:HB3	2.02	0.41
1:C:747:LEU:HD12	1:C:786:VAL:HB	2.01	0.41
1:C:905:TRP:HB2	1:C:947:MET:HE3	2.02	0.41
1:D:805:HIS:O	1:D:809:ILE:HG13	2.20	0.41
1:B:731:TRP:CZ3	1:B:733:PRO:HG3	2.56	0.41
1:D:766:MET:HE3	1:D:777:LEU:HD22	2.02	0.41
1:D:986:ARG:H	1:D:986:ARG:HG3	1.67	0.41
1:A:762:GLU:CD	1:A:860:LYS:HB2	2.41	0.41
1:B:908[B]:MET:SD	1:B:943:VAL:HG11	2.60	0.41
1:C:723:PHE:CZ	1:C:745:LYS:HD2	2.56	0.41
1:D:900:TYR:O	1:D:904:VAL:HG23	2.21	0.40
1:D:860:LYS:HB3	1:D:860:LYS:HE3	1.77	0.40
1:C:793:MET:CE	1:C:852:LYS:HD3	2.51	0.40
1:B:962:ARG:HH11	1:B:962:ARG:HD2	1.74	0.40
1:A:705:ARG:HE	1:A:707:LEU:HD23	1.87	0.40
1:D:745:LYS:HE3	1:D:745:LYS:HB2	1.56	0.40
1:C:990:PRO:HB2	1:C:994:ASP:HB2	2.04	0.40
1:B:731:TRP:CH2	1:B:733:PRO:HG3	2.56	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	289/331 (87%)	280 (97%)	8 (3%)	1 (0%)	41	55
1	B	280/331 (85%)	270 (96%)	9 (3%)	1 (0%)	34	48
1	C	287/331 (87%)	276 (96%)	10 (4%)	1 (0%)	41	55
1	D	289/331 (87%)	279 (96%)	9 (3%)	1 (0%)	41	55
All	All	1145/1324 (86%)	1105 (96%)	36 (3%)	4 (0%)	34	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	859	ALA
1	C	858	LEU
1	B	858	LEU
1	A	1011	VAL

### 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	264/290 (91%)	257 (97%)	7 (3%)	44	65
1	B	255/290 (88%)	249 (98%)	6 (2%)	49	68
1	C	263/290 (91%)	252 (96%)	11 (4%)	30	47
1	D	262/290 (90%)	252 (96%)	10 (4%)	33	51
All	All	1044/1160 (90%)	1010 (97%)	34 (3%)	38	57

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	723	PHE
1	D	754	LYS
1	D	757	LYS
1	D	783	THR
1	D	784	SER
1	D	860	LYS
1	D	960	LYS
1	D	986	ARG
1	D	998	TYR
1	D	1007	MET
1	A	745	LYS
1	A	748	ARG
1	A	783	THR
1	A	945	MET
1	A	1006	ASP
1	A	1011	VAL
1	A	1012	ASP
1	C	748	ARG
1	C	754	LYS
1	C	760	LEU
1	C	783	THR
1	C	784	SER
1	C	942	ASP
1	C	949	LYS
1	C	960	LYS
1	C	962	ARG
1	C	998	TYR
1	C	1006	ASP
1	B	808	ASN
1	B	860	LYS
1	B	918	ILE
1	B	931	GLU
1	B	960	LYS
1	B	962	ARG

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

Mol	Chain	Res	Type
1	A	816	ASN
1	A	976	GLN
1	C	701	GLN

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Mol	Chain	Res	Type
1	C	773	HIS
1	B	808	ASN

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

4 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	VO7	A	1101	1	38,39,39	2.48	13 (34%)	50,54,54	3.35	19 (38%)
2	VO7	C	1101	1	38,39,39	2.21	11 (28%)	50,54,54	3.33	15 (30%)
2	VO7	B	1101	1	38,39,39	2.40	17 (44%)	50,54,54	3.24	19 (38%)
2	VO7	D	1101	1	38,39,39	2.21	15 (39%)	50,54,54	3.56	14 (28%)

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	VO7	A	1101	1	-	4/21/25/25	0/4/4/4
2	VO7	C	1101	1	-	6/21/25/25	0/4/4/4
2	VO7	B	1101	1	-	6/21/25/25	0/4/4/4
2	VO7	D	1101	1	-	8/21/25/25	0/4/4/4

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1101	VO7	C16-C15	6.60	1.59	1.48
2	D	1101	VO7	C15-N14	5.54	1.45	1.35
2	A	1101	VO7	C13-C05	5.39	1.46	1.40
2	B	1101	VO7	C20-N21	5.22	1.54	1.39
2	A	1101	VO7	C20-N21	5.19	1.54	1.39
2	D	1101	VO7	C16-C15	5.17	1.56	1.48
2	A	1101	VO7	C05-N06	5.02	1.52	1.41
2	A	1101	VO7	C16-C15	4.86	1.56	1.48
2	A	1101	VO7	C19-C13	4.73	1.46	1.39
2	B	1101	VO7	C16-C15	4.72	1.56	1.48
2	B	1101	VO7	C09-N10	4.56	1.62	1.46
2	A	1101	VO7	C15-N14	4.53	1.43	1.35
2	C	1101	VO7	C15-N14	4.42	1.43	1.35
2	C	1101	VO7	C22-N36	4.39	1.40	1.34
2	D	1101	VO7	C13-N14	4.38	1.50	1.41
2	C	1101	VO7	C13-N14	4.32	1.50	1.41
2	A	1101	VO7	C22-N36	4.29	1.40	1.34
2	C	1101	VO7	C26-N27	3.68	1.44	1.36
2	B	1101	VO7	C08-N06	3.64	1.58	1.46
2	B	1101	VO7	C12-N10	3.62	1.57	1.46
2	A	1101	VO7	C26-N27	3.58	1.44	1.36
2	B	1101	VO7	C22-N36	3.51	1.39	1.34
2	C	1101	VO7	C20-N21	3.34	1.49	1.39
2	B	1101	VO7	C22-N21	3.18	1.43	1.36
2	B	1101	VO7	C07-N06	3.17	1.51	1.46
2	B	1101	VO7	C15-N14	3.16	1.41	1.35
2	D	1101	VO7	C26-N27	3.12	1.43	1.36
2	D	1101	VO7	C22-N36	3.10	1.39	1.34
2	D	1101	VO7	C20-N21	3.08	1.48	1.39
2	B	1101	VO7	C11-N10	3.05	1.55	1.46
2	C	1101	VO7	C11-N10	3.01	1.55	1.46
2	D	1101	VO7	C09-N10	2.91	1.56	1.46
2	B	1101	VO7	C19-C13	2.88	1.44	1.39
2	A	1101	VO7	C04-C03	2.87	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	VO7	C09-C08	2.84	1.62	1.51
2	A	1101	VO7	C22-N21	2.84	1.42	1.36
2	B	1101	VO7	C04-C03	2.76	1.43	1.38
2	B	1101	VO7	C13-C05	2.74	1.43	1.40
2	B	1101	VO7	C05-N06	2.71	1.47	1.41
2	C	1101	VO7	C13-C05	2.70	1.43	1.40
2	C	1101	VO7	C08-N06	2.65	1.55	1.46
2	D	1101	VO7	C11-N10	2.62	1.54	1.46
2	C	1101	VO7	C09-N10	2.58	1.55	1.46
2	A	1101	VO7	C13-N14	2.52	1.46	1.41
2	D	1101	VO7	O02-C03	2.47	1.41	1.37
2	B	1101	VO7	C26-N27	2.44	1.41	1.36
2	D	1101	VO7	C19-C13	2.41	1.43	1.39
2	D	1101	VO7	C22-N21	2.37	1.41	1.36
2	A	1101	VO7	C25-C24	-2.36	1.46	1.49
2	D	1101	VO7	C04-C03	2.35	1.43	1.38
2	A	1101	VO7	C12-N10	2.32	1.53	1.46
2	C	1101	VO7	C09-C08	2.27	1.60	1.51
2	D	1101	VO7	C20-C03	2.26	1.45	1.40
2	D	1101	VO7	C13-C05	2.18	1.43	1.40
2	B	1101	VO7	O02-C03	2.15	1.40	1.37
2	D	1101	VO7	C04-C05	2.09	1.43	1.39

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1101	VO7	C16-C15-N14	21.05	127.71	113.84
2	C	1101	VO7	C16-C15-N14	19.02	126.37	113.84
2	A	1101	VO7	C16-C15-N14	18.21	125.84	113.84
2	B	1101	VO7	C16-C15-N14	17.12	125.12	113.84
2	B	1101	VO7	C09-C08-N06	5.39	120.91	112.31
2	D	1101	VO7	O18-C15-C16	-5.23	114.37	122.72
2	B	1101	VO7	N36-C22-N23	-5.19	121.63	126.55
2	A	1101	VO7	N36-C22-N23	-4.73	122.06	126.55
2	B	1101	VO7	O18-C15-N14	-4.63	117.74	123.05
2	A	1101	VO7	C19-C20-C03	-4.58	113.38	118.91
2	D	1101	VO7	N36-C22-N23	-4.58	122.21	126.55
2	C	1101	VO7	O18-C15-N14	-4.54	117.85	123.05
2	D	1101	VO7	O18-C15-N14	-4.50	117.89	123.05
2	A	1101	VO7	C05-C04-C03	4.49	127.48	117.91
2	A	1101	VO7	O18-C15-C16	-4.45	115.61	122.72
2	C	1101	VO7	N36-C22-N23	-4.37	122.40	126.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1101	VO7	O18-C15-C16	-4.34	115.78	122.72
2	A	1101	VO7	C04-C05-C13	-4.32	113.66	120.04
2	C	1101	VO7	C09-C08-N06	4.24	119.06	112.31
2	B	1101	VO7	C05-C04-C03	4.10	126.64	117.91
2	A	1101	VO7	C24-N23-C22	4.10	120.06	116.69
2	B	1101	VO7	C24-N23-C22	3.99	119.98	116.69
2	D	1101	VO7	O02-C03-C20	3.95	119.65	114.80
2	A	1101	VO7	O18-C15-N14	-3.92	118.56	123.05
2	B	1101	VO7	C19-C20-C03	-3.89	114.22	118.91
2	C	1101	VO7	C24-N23-C22	3.88	119.89	116.69
2	C	1101	VO7	C05-C04-C03	3.77	125.94	117.91
2	A	1101	VO7	C24-C25-C29	3.66	131.05	123.90
2	B	1101	VO7	C24-C25-C29	3.63	131.00	123.90
2	D	1101	VO7	C05-C04-C03	3.58	125.54	117.91
2	B	1101	VO7	O18-C15-C16	-3.54	117.07	122.72
2	D	1101	VO7	C24-N23-C22	3.43	119.51	116.69
2	C	1101	VO7	C19-C20-C03	-3.23	115.00	118.91
2	D	1101	VO7	C19-C20-C03	-3.20	115.05	118.91
2	B	1101	VO7	C04-C05-C13	-3.14	115.41	120.04
2	D	1101	VO7	C19-C13-C05	-3.12	116.39	119.67
2	A	1101	VO7	C25-C24-N23	2.91	119.95	116.14
2	C	1101	VO7	C19-C13-C05	-2.84	116.67	119.67
2	A	1101	VO7	O02-C03-C20	2.75	118.17	114.80
2	A	1101	VO7	C08-N06-C05	2.74	128.25	117.16
2	C	1101	VO7	C30-C29-C28	-2.69	114.60	118.17
2	B	1101	VO7	C30-C29-C28	-2.57	114.76	118.17
2	C	1101	VO7	O02-C03-C20	2.56	117.94	114.80
2	A	1101	VO7	C04-C05-N06	-2.56	117.48	120.93
2	C	1101	VO7	C04-C05-C13	-2.56	116.26	120.04
2	A	1101	VO7	C30-C29-C28	-2.54	114.79	118.17
2	C	1101	VO7	C24-C25-C29	2.54	128.86	123.90
2	D	1101	VO7	C24-C25-C29	2.45	128.69	123.90
2	B	1101	VO7	C19-C13-C05	-2.35	117.19	119.67
2	C	1101	VO7	C07-N06-C05	2.30	122.23	115.89
2	D	1101	VO7	C30-C29-C28	-2.27	115.16	118.17
2	D	1101	VO7	C13-C19-C20	2.26	127.94	121.59
2	B	1101	VO7	C31-C30-C29	2.24	123.99	120.89
2	D	1101	VO7	C04-C03-C20	-2.21	117.88	120.54
2	B	1101	VO7	C25-C24-N23	2.19	119.02	116.14
2	B	1101	VO7	C01-O02-C03	2.17	120.80	117.53
2	A	1101	VO7	C13-C19-C20	2.16	127.67	121.59
2	A	1101	VO7	C19-C13-C05	-2.15	117.41	119.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	VO7	C19-C20-N21	2.13	125.47	121.05
2	B	1101	VO7	C13-C19-C20	2.11	127.53	121.59
2	A	1101	VO7	C17-C16-C15	2.10	124.67	122.27
2	B	1101	VO7	O02-C03-C20	2.09	117.36	114.80
2	B	1101	VO7	C13-N14-C15	2.06	131.91	127.15
2	C	1101	VO7	C13-C19-C20	2.06	127.38	121.59
2	B	1101	VO7	C19-C20-N21	2.03	125.27	121.05
2	A	1101	VO7	C31-C30-C29	2.01	123.68	120.89
2	D	1101	VO7	C03-C20-N21	2.00	121.97	117.78

There are no chirality outliers.

All (24) torsion outliers are listed below:

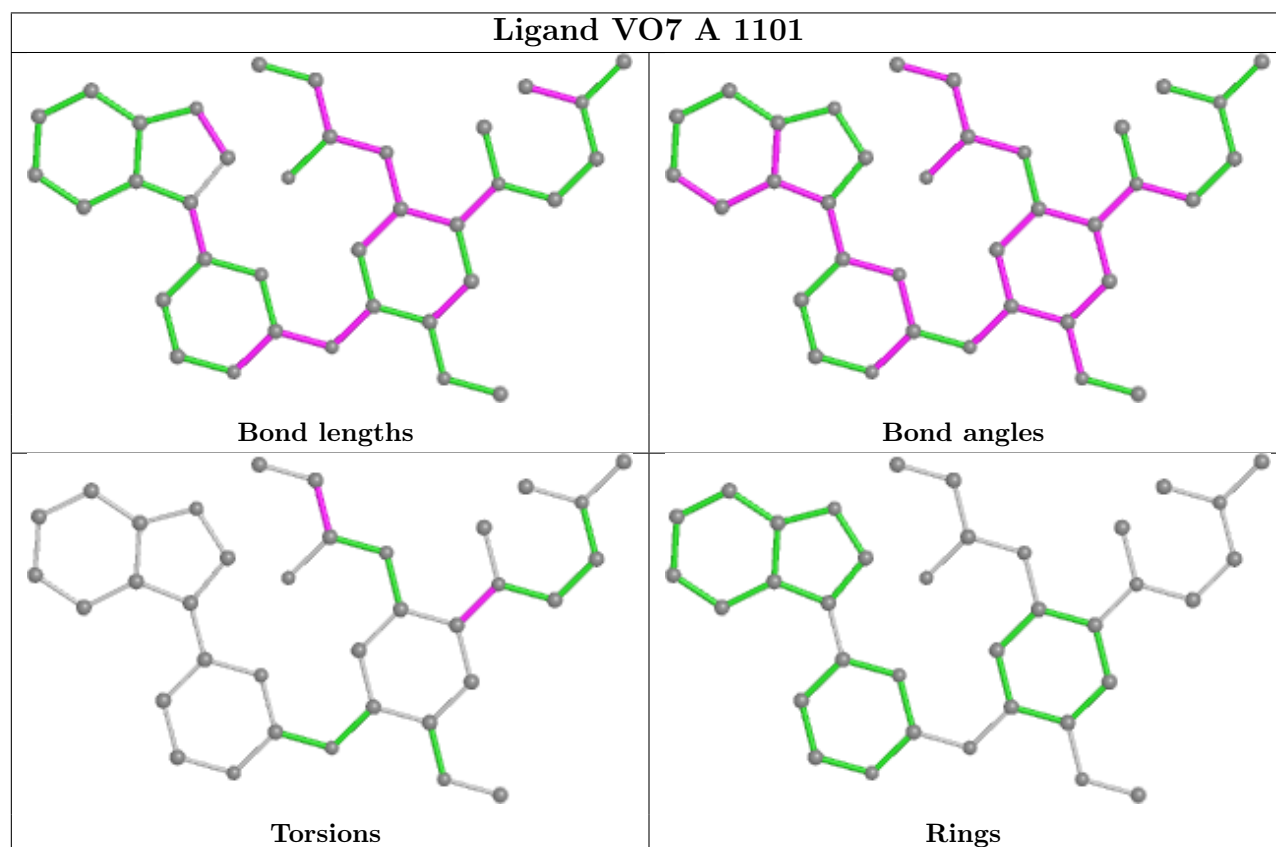
Mol	Chain	Res	Type	Atoms
2	D	1101	VO7	N14-C15-C16-C17
2	D	1101	VO7	O18-C15-C16-C17
2	D	1101	VO7	C09-C08-N06-C07
2	A	1101	VO7	N14-C15-C16-C17
2	A	1101	VO7	O18-C15-C16-C17
2	C	1101	VO7	N14-C15-C16-C17
2	C	1101	VO7	O18-C15-C16-C17
2	C	1101	VO7	C09-C08-N06-C07
2	B	1101	VO7	N14-C15-C16-C17
2	B	1101	VO7	O18-C15-C16-C17
2	B	1101	VO7	C09-C08-N06-C07
2	D	1101	VO7	C08-C09-N10-C11
2	C	1101	VO7	C08-C09-N10-C11
2	C	1101	VO7	N06-C08-C09-N10
2	D	1101	VO7	N06-C08-C09-N10
2	B	1101	VO7	C13-C05-N06-C08
2	D	1101	VO7	C13-C05-N06-C07
2	B	1101	VO7	C09-C08-N06-C05
2	A	1101	VO7	C13-C05-N06-C08
2	D	1101	VO7	C04-C05-N06-C07
2	A	1101	VO7	C04-C05-N06-C08
2	D	1101	VO7	C09-C08-N06-C05
2	C	1101	VO7	C09-C08-N06-C05
2	B	1101	VO7	C04-C05-N06-C08

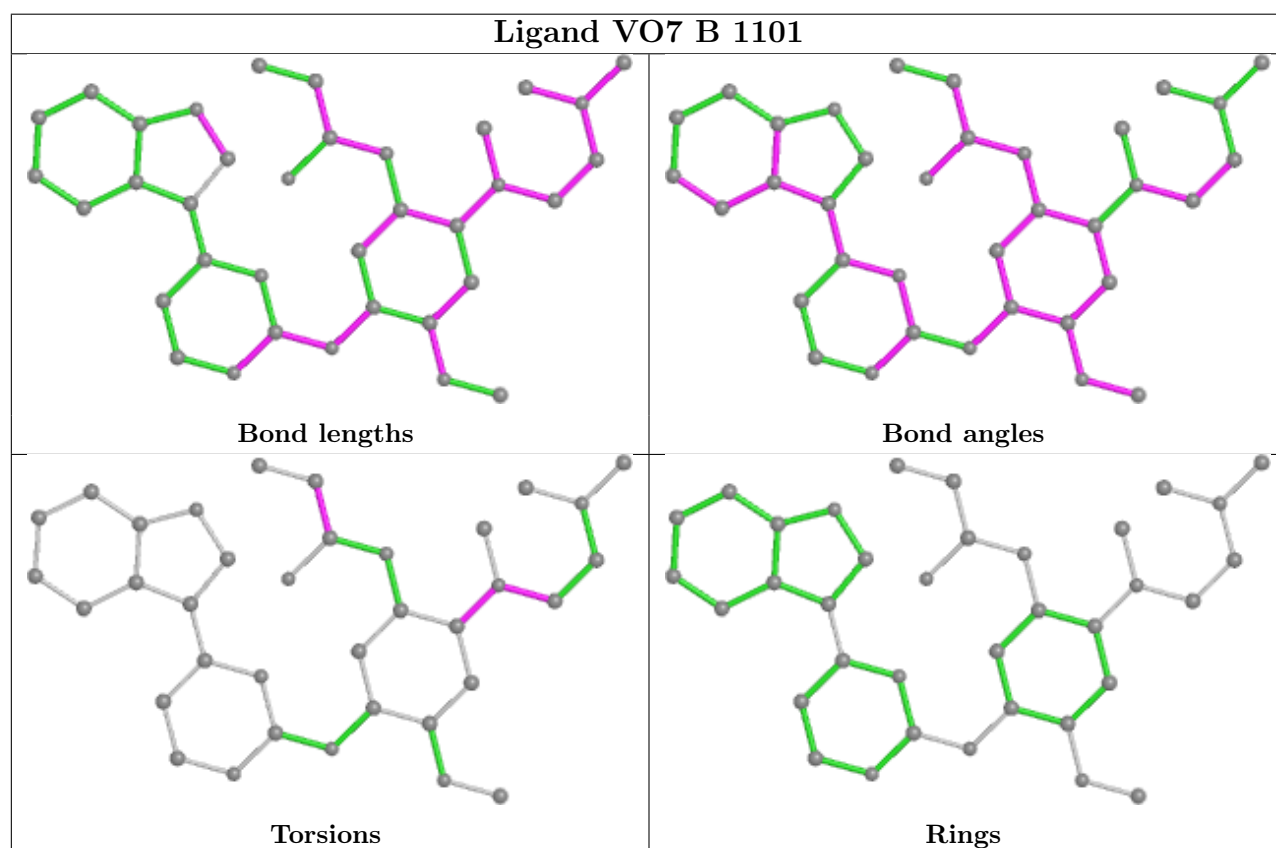
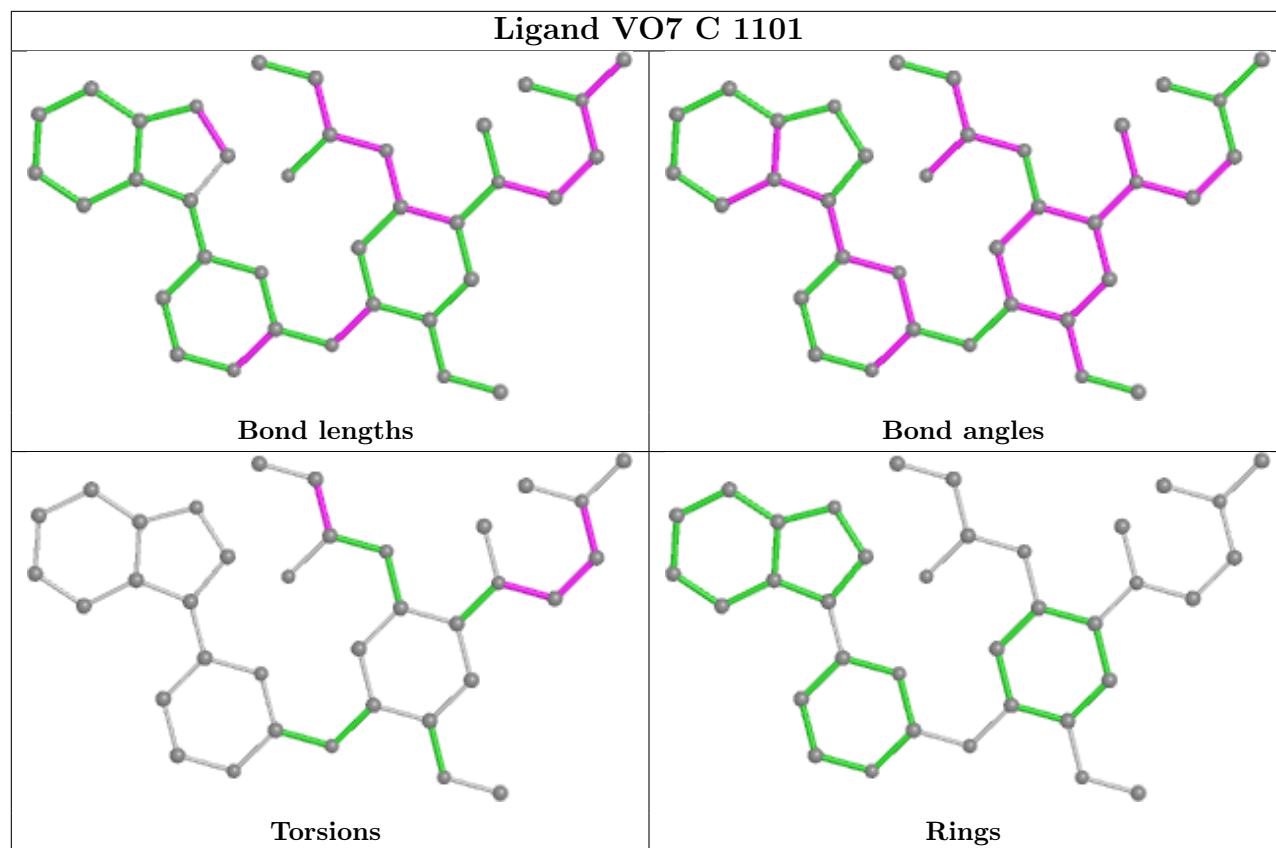
There are no ring outliers.

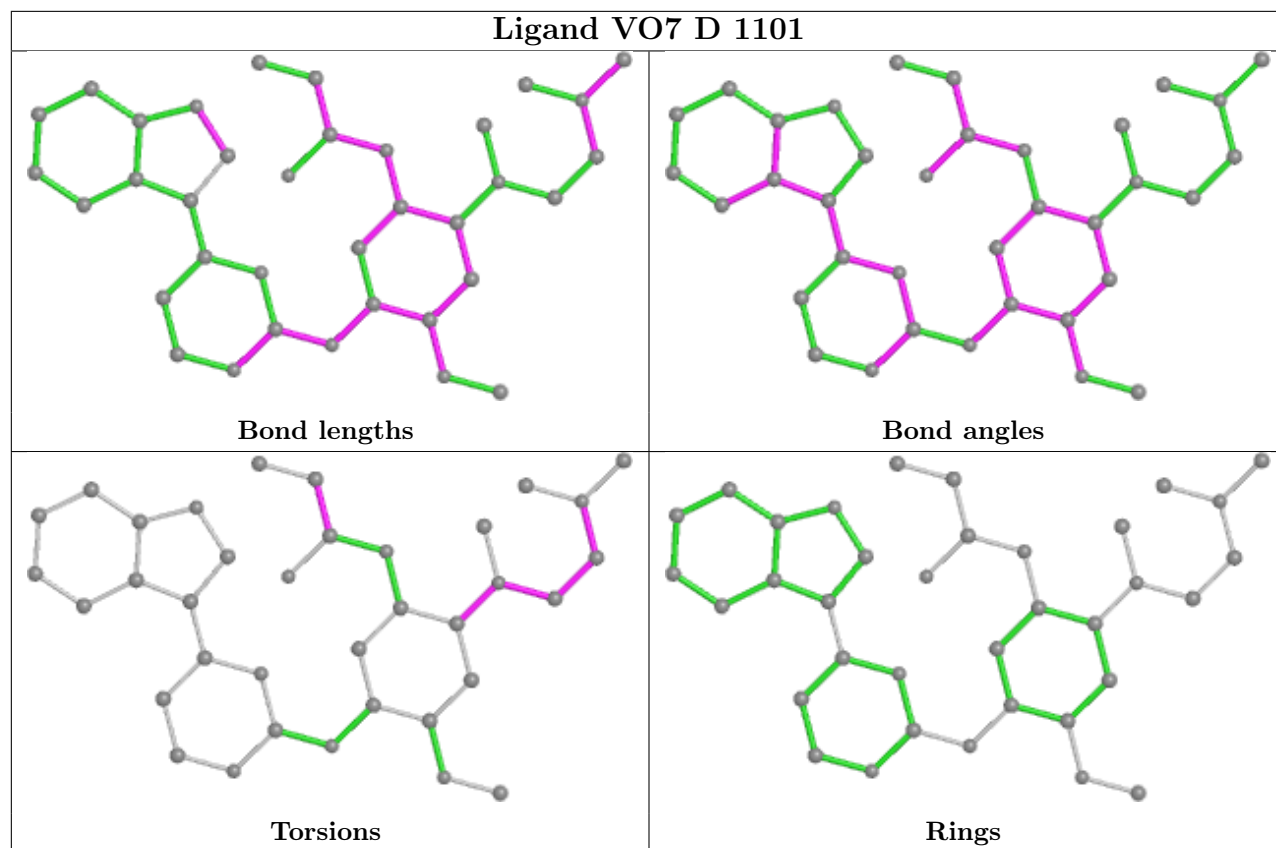
4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	VO7	1	0
2	C	1101	VO7	3	0
2	B	1101	VO7	1	0
2	D	1101	VO7	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 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	294/331 (88%)	0.05	8 (2%) 54 52	30, 46, 69, 81	0
1	B	285/331 (86%)	0.18	9 (3%) 47 46	30, 48, 80, 93	0
1	C	292/331 (88%)	0.25	9 (3%) 49 47	35, 54, 85, 103	0
1	D	292/331 (88%)	0.08	7 (2%) 59 57	31, 44, 71, 86	0
All	All	1163/1324 (87%)	0.14	33 (2%) 53 51	30, 48, 78, 103	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	723	PHE	7.3
1	D	988	HIS	5.7
1	C	860	LYS	5.0
1	B	784	SER	4.3
1	C	859	ALA	4.2
1	D	722	ALA	3.9
1	D	858	LEU	3.9
1	C	986	ARG	3.8
1	C	988	HIS	3.4
1	D	748	ARG	3.4
1	C	977	ARG	3.2
1	B	752	SER	3.2
1	D	857	GLY	3.1
1	A	988	HIS	3.0
1	C	857	GLY	2.9
1	B	981	ILE	2.9
1	B	751	THR	2.8
1	B	992	PRO	2.7
1	A	808	ASN	2.6
1	C	929	LYS	2.6
1	A	860	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	917	GLY	2.5
1	D	700	ASN	2.4
1	A	737	LYS	2.4
1	C	753	PRO	2.4
1	B	708	LYS	2.2
1	A	748	ARG	2.2
1	A	1010	VAL	2.1
1	C	992	PRO	2.1
1	B	734	GLU	2.1
1	B	921	SER	2.0
1	A	919	PRO	2.0
1	B	860	LYS	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.

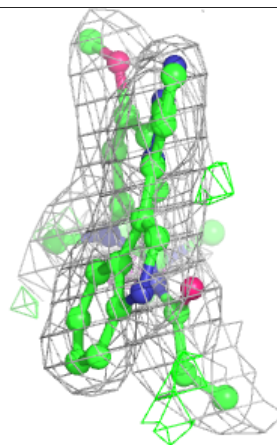
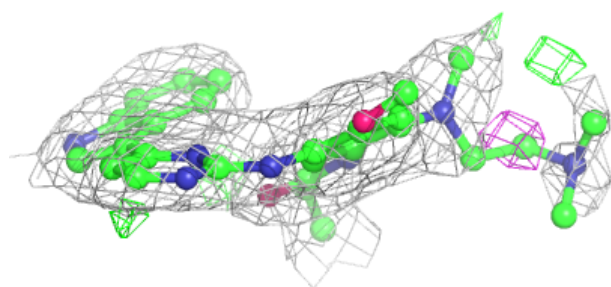
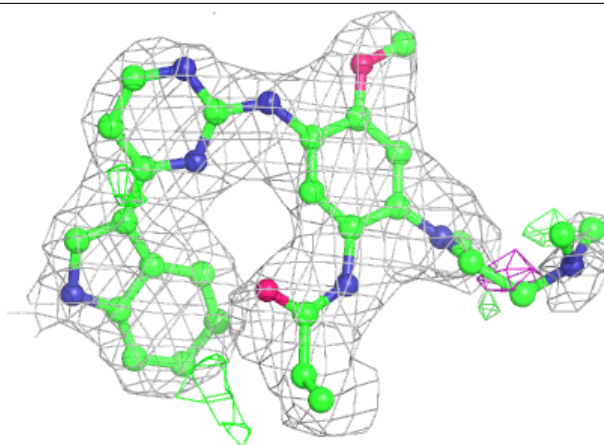
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	VO7	C	1101	36/36	0.91	0.19	39,45,69,77	0
2	VO7	A	1101	36/36	0.92	0.20	39,50,67,72	0
2	VO7	D	1101	36/36	0.92	0.17	39,47,69,74	0
2	VO7	B	1101	36/36	0.92	0.21	37,47,66,71	0

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.



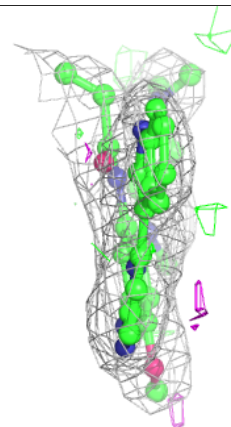
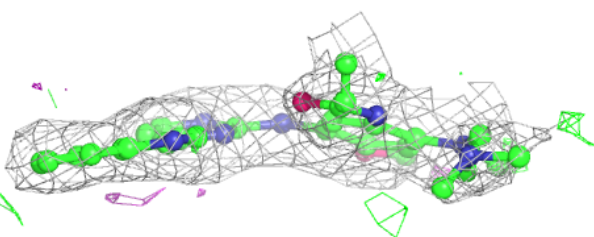
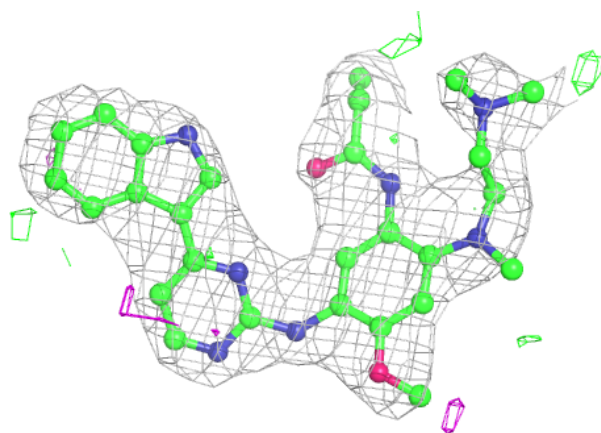
**Electron density around VO7 C 1101:**

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

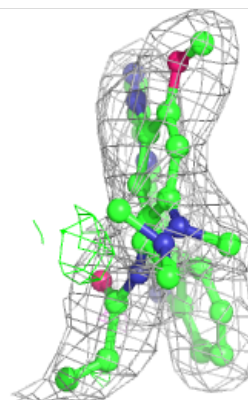
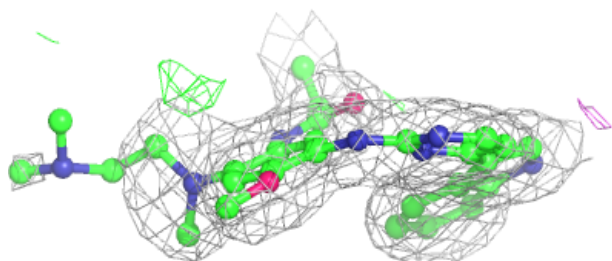
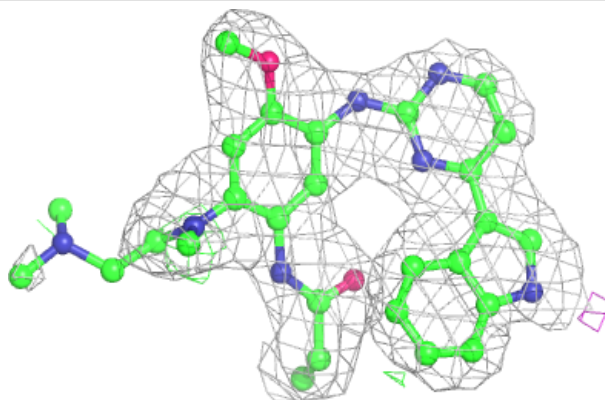


**Electron density around VO7 A 1101:**

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

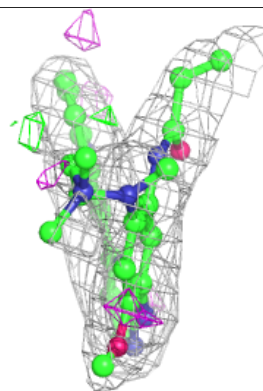
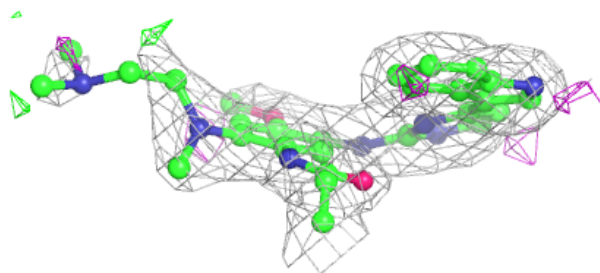
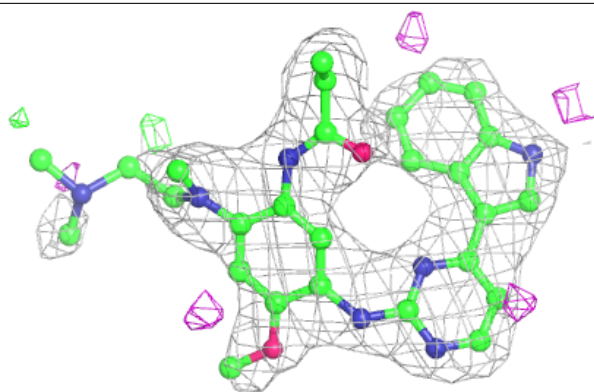
**Electron density around VO7 D 1101:**

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



**Electron density around VO7 B 1101:**

$2mF_o - DF_c$  (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.