



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 04:53 PM BST

PDB ID : 5XIP  
Title : Crystal Structure of Eimeria tenella Prolyl-tRNA Synthetase (EtPRS) in complex with Halofuginone  
Authors : Jain, V.; Manickam, Y.; Sharma, A.  
Deposited on : 2017-04-26  
Resolution : 3.10 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

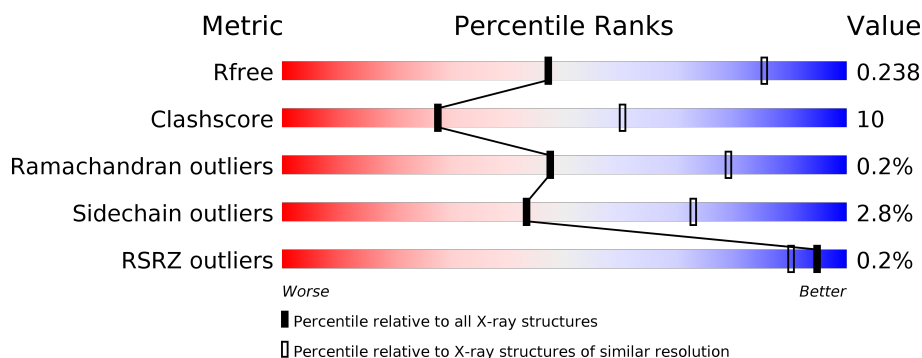
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	498	<div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div>
1	B	498	<div> <div>77%</div> <div>18%</div> <div>• 5%</div> </div>
1	C	498	<div> <div>71%</div> <div>21%</div> <div>• 7%</div> </div>
1	D	498	<div> <div>65%</div> <div>25%</div> <div>• 8%</div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 15058 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 Prolyl-tRNA synthetase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	0	0	0
			3767	2420	652	675	20			
1	B	475	Total	C	N	O	S	0	1	0
			3790	2430	653	688	19			
1	C	463	Total	C	N	O	S	0	1	0
			3633	2337	625	652	19			
1	D	457	Total	C	N	O	S	0	2	0
			3628	2334	622	652	20			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	250	GLY	-	expression tag	UNP U6KWI1
A	251	ALA	-	expression tag	UNP U6KWI1
A	252	MET	-	expression tag	UNP U6KWI1
B	250	GLY	-	expression tag	UNP U6KWI1
B	251	ALA	-	expression tag	UNP U6KWI1
B	252	MET	-	expression tag	UNP U6KWI1
C	250	GLY	-	expression tag	UNP U6KWI1
C	251	ALA	-	expression tag	UNP U6KWI1
C	252	MET	-	expression tag	UNP U6KWI1
D	250	GLY	-	expression tag	UNP U6KWI1
D	251	ALA	-	expression tag	UNP U6KWI1
D	252	MET	-	expression tag	UNP U6KWI1

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>12</sub>P<sub>3</sub>).

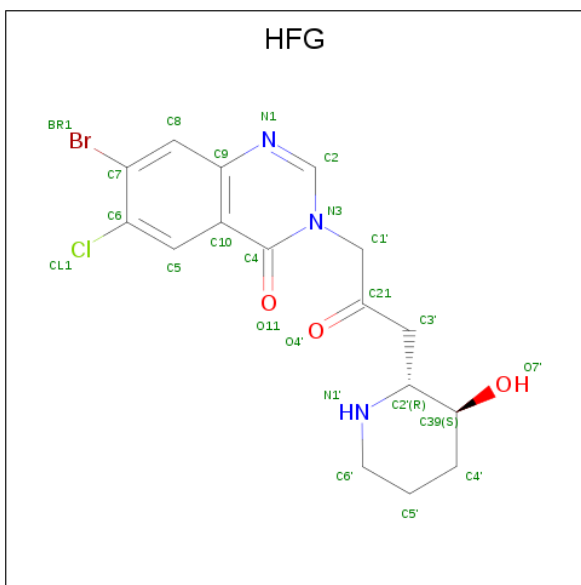


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	B	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	C	1	Total 31	C 10	N 6	O 12	P 3	0	0
2	D	1	Total 31	C 10	N 6	O 12	P 3	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is 7-bromo-6-chloro-3-[3-[(2R,3S)-3-hydroxypiperidin-2-yl]-2-oxopropyl]quinazolin-4(3H)-one (three-letter code: HFG) (formula: C<sub>16</sub>H<sub>17</sub>BrClN<sub>3</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		
4	B	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		
4	C	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		
4	D	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		

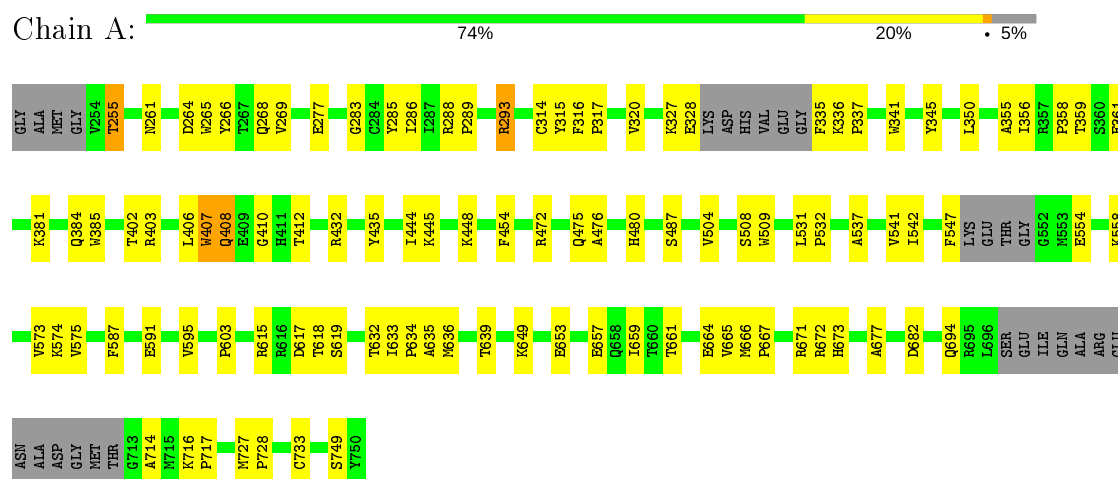
- Molecule 5 is water.

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

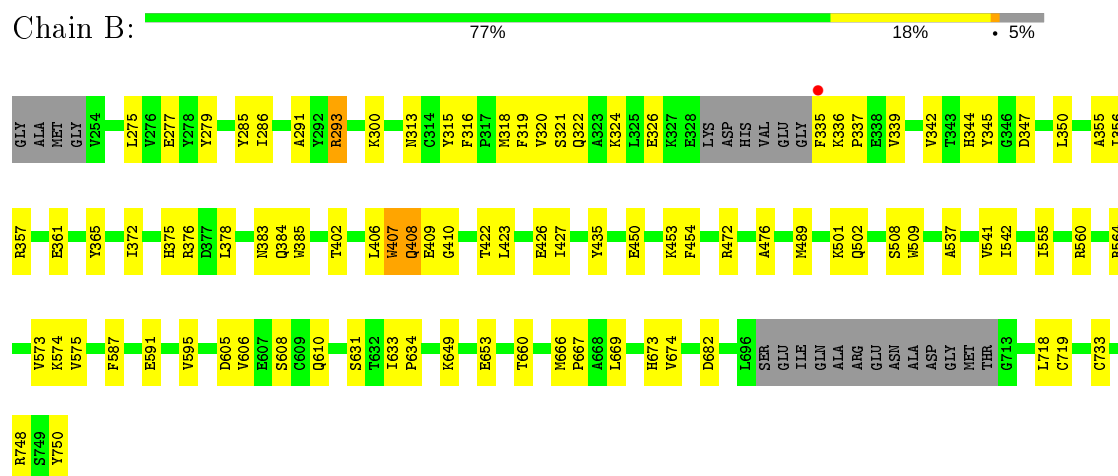
### 3 Residue-property plots

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: Prolyl-tRNA synthetase, putative

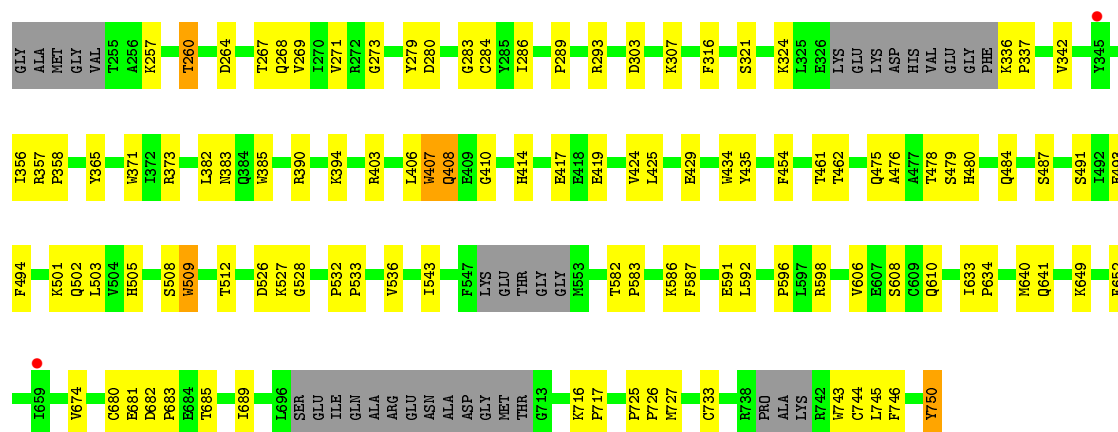


- Molecule 1: Prolyl-tRNA synthetase, putative



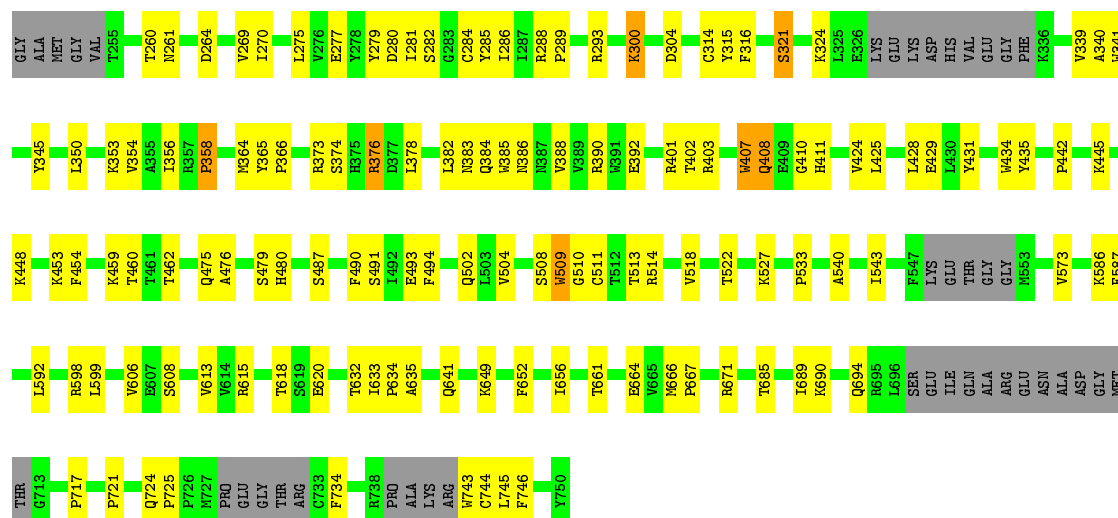
- Molecule 1: Prolyl-tRNA synthetase, putative





• Molecule 1: Prolyl-tRNA synthetase, putative

Chain D: 65% 25% 8%



## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.76Å 138.76Å 425.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.77 – 3.10 41.77 – 3.10	Depositor EDS
% Data completeness (in resolution range)	95.5 (41.77-3.10) 95.6 (41.77-3.10)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.31 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.180 , 0.230 0.193 , 0.238	Depositor DCC
$R_{free}$ test set	2003 reflections (3.78%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.6	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , -8.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.459 for -h-k,k,-l	Xtriage
Reported twinning fraction	0.588 for H, K, L 0.412 for K, H, -L	Depositor
Outliers	0 of 53009 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	15058	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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: MG, ANP, HFG

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	0.71	0/3865	0.81	0/5248
1	B	0.71	0/3890	0.82	0/5283
1	C	0.65	0/3727	0.80	0/5065
1	D	0.65	0/3726	0.82	0/5062
All	All	0.68	0/15208	0.81	0/20658

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	3767	0	3694	69	0
1	B	3790	0	3698	65	0
1	C	3633	0	3524	73	0
1	D	3628	0	3514	99	0
2	A	31	0	13	2	0
2	B	31	0	13	0	0
2	C	31	0	13	2	0
2	D	31	0	13	4	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	24	0	17	6	0
4	B	24	0	17	7	0
4	C	24	0	17	6	0
4	D	24	0	17	4	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
All	All	15058	0	14550	296	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 10.

All (296) 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:B:407:TRP:CH2	4:B:1003:HFG:H8	1.92	1.04
1:A:407:TRP:CH2	4:A:1003:HFG:H8	2.01	0.95
1:A:356:ILE:HD11	1:B:356:ILE:HD11	1.56	0.88
1:A:407:TRP:CZ3	4:A:1003:HFG:H8	2.08	0.88
1:A:454:PHE:CE1	4:A:1003:HFG:H15	2.13	0.83
1:A:345:TYR:HD1	1:A:350:LEU:HD23	1.51	0.75
1:D:279:TYR:HB3	1:D:284:CYS:O	1.86	0.74
1:D:390:ARG:NH1	2:D:1001:ANP:O2B	2.20	0.74
1:A:359:THR:OG1	1:A:361:GLU:OE1	2.03	0.74
1:B:407:TRP:HH2	4:B:1003:HFG:H8	1.52	0.72
1:D:429:GLU:HA	1:D:429:GLU:OE1	1.89	0.71
1:A:293:ARG:HD3	1:A:537:ALA:HA	1.73	0.71
1:A:435:TYR:CZ	1:A:476:ALA:HB1	2.25	0.71
1:D:435:TYR:CZ	1:D:476:ALA:HB1	2.25	0.71
1:B:454:PHE:CZ	4:B:1003:HFG:H1	2.26	0.70
1:D:378:LEU:HD11	1:D:504:VAL:HG12	1.73	0.70
1:D:725:PRO:HG2	1:D:743:TRP:CD1	2.26	0.70
4:D:1003:HFG:O4'	4:D:1003:HFG:N1'	2.26	0.68
1:A:385:TRP:CZ3	1:A:408:GLN:HB2	2.28	0.68
1:C:487:SER:O	1:C:491:SER:N	2.26	0.68
1:C:414:HIS:HD1	1:C:419:GLU:HG2	1.59	0.67
1:D:480:HIS:NE2	4:D:1003:HFG:O7'	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:435:TYR:CZ	1:B:476:ALA:HB1	2.29	0.67
1:B:407:TRP:CZ3	4:B:1003:HFG:H8	2.30	0.67
1:A:435:TYR:CE2	1:A:476:ALA:HB1	2.29	0.66
1:C:484:GLN:O	1:C:487:SER:HB2	1.96	0.66
1:A:649:LYS:O	1:A:653:GLU:HG2	1.95	0.66
1:B:372:ILE:HD13	1:B:378:LEU:HD21	1.77	0.65
1:B:339:VAL:HG22	4:B:1003:HFG:BR1	2.51	0.65
1:D:744:CYS:SG	1:D:745:LEU:N	2.70	0.65
1:B:320:VAL:HG12	1:B:355:ALA:HB3	1.79	0.65
1:C:407:TRP:CH2	4:C:1003:HFG:H8	2.31	0.65
1:D:618:THR:OG1	1:D:620:GLU:HG2	1.98	0.64
1:A:406:LEU:HD21	1:B:315:TYR:CE2	2.34	0.62
1:C:596:PRO:HB3	1:C:640:MET:SD	2.40	0.62
1:D:724:GLN:OE1	1:D:743:TRP:HB2	1.98	0.62
1:B:372:ILE:HD13	1:B:378:LEU:CD2	2.31	0.60
1:B:293:ARG:CG	1:B:293:ARG:HH11	2.14	0.60
1:A:283:GLY:HA3	1:A:403:ARG:HG3	1.85	0.59
1:D:285:TYR:HE1	1:D:402:THR:HG22	1.67	0.58
1:C:280:ASP:OD2	1:D:324:LYS:HD2	2.03	0.58
1:B:385:TRP:CZ3	1:B:408:GLN:HB2	2.39	0.58
1:C:633:ILE:HB	1:C:634:PRO:HD3	1.84	0.58
1:B:422:THR:O	1:B:426:GLU:HG2	2.04	0.58
1:B:285:TYR:HE1	1:B:402:THR:HG22	1.69	0.57
1:C:744:CYS:SG	1:C:745:LEU:N	2.78	0.57
1:A:406:LEU:HD21	1:B:315:TYR:CZ	2.39	0.57
1:C:424:VAL:HG12	1:C:425:LEU:HD23	1.87	0.57
1:A:618:THR:O	1:A:619:SER:HB2	2.05	0.57
1:C:501:LYS:O	1:C:502:GLN:HG2	2.04	0.56
1:A:345:TYR:CD1	1:A:350:LEU:HD23	2.38	0.56
1:D:339:VAL:HG12	1:D:340:ALA:O	2.06	0.56
1:A:385:TRP:CE3	1:A:408:GLN:HB2	2.41	0.56
1:D:649:LYS:O	1:D:652:PHE:HB3	2.05	0.56
1:B:277:GLU:HB3	1:B:286:ILE:HB	1.88	0.56
1:B:472:ARG:HH12	1:B:748:ARG:HB3	1.71	0.56
1:C:725:PRO:O	1:C:743:TRP:NE1	2.38	0.56
1:B:321:SER:OG	1:B:324:LYS:HB2	2.07	0.55
1:C:435:TYR:CZ	1:C:476:ALA:HB1	2.41	0.55
1:A:657:GLU:OE2	1:A:673:HIS:ND1	2.39	0.55
1:B:560:ARG:O	1:B:564:ARG:HG3	2.06	0.55
1:B:345:TYR:HD1	1:B:350:LEU:HD23	1.72	0.55
1:B:649:LYS:O	1:B:653:GLU:HG2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:394:LYS:NZ	1:C:750:TYR:OXT	2.40	0.55
1:A:381:LYS:HA	1:A:412:THR:HG22	1.89	0.55
1:A:694:GLN:HB2	1:A:714:ALA:HB2	1.87	0.55
1:C:279:TYR:HB3	1:C:284:CYS:O	2.07	0.55
2:D:1001:ANP:O1A	4:D:1003:HFG:H2	2.08	0.55
1:D:383:ASN:OD1	1:D:408:GLN:HG2	2.07	0.55
1:A:487:SER:OG	1:A:504:VAL:HG23	2.06	0.54
1:C:480:HIS:NE2	4:C:1003:HFG:O7'	2.32	0.54
2:A:1001:ANP:O1A	4:A:1003:HFG:H17	2.07	0.54
1:C:407:TRP:HB3	1:C:512:THR:HG22	1.90	0.54
1:A:315:TYR:CZ	1:B:406:LEU:HD21	2.42	0.54
1:D:633:ILE:HB	1:D:634:PRO:HD3	1.89	0.54
1:B:407:TRP:CH2	4:B:1003:HFG:C5'	2.80	0.53
1:C:390:ARG:NH1	2:C:1001:ANP:O1A	2.40	0.53
1:B:587:PHE:O	1:B:591:GLU:HG3	2.09	0.53
1:C:371:TRP:NE1	1:D:277:GLU:OE2	2.40	0.53
1:B:682:ASP:OD1	1:B:682:ASP:C	2.47	0.53
1:D:316:PHE:CE1	1:D:364:MET:HG2	2.44	0.53
1:C:286:ILE:HD13	1:D:316:PHE:CE2	2.44	0.53
1:B:293:ARG:HH11	1:B:293:ARG:HG2	1.73	0.53
1:B:320:VAL:CG1	1:B:355:ALA:HB3	2.39	0.53
1:A:327:LYS:O	1:A:328:GLU:CB	2.56	0.53
1:B:336:LYS:N	1:B:337:PRO:CD	2.72	0.53
1:C:484:GLN:HG2	1:C:503:LEU:HB3	1.90	0.52
1:D:390:ARG:NH1	2:D:1001:ANP:O2A	2.36	0.52
1:A:587:PHE:O	1:A:591:GLU:HG3	2.10	0.52
1:B:435:TYR:CE2	1:B:476:ALA:HB1	2.44	0.52
1:A:573:VAL:HG12	1:A:574:LYS:N	2.25	0.52
1:C:303:ASP:OD1	1:C:307:LYS:HD2	2.10	0.52
1:B:385:TRP:CE3	1:B:408:GLN:HB2	2.45	0.52
1:B:669:LEU:HB3	1:B:748:ARG:NH1	2.24	0.52
1:B:472:ARG:NH1	1:B:748:ARG:HB3	2.25	0.52
1:A:480:HIS:NE2	4:A:1003:HFG:O7'	2.31	0.52
1:D:533:PRO:HD2	1:D:641:GLN:OE1	2.10	0.51
1:C:257:LYS:HA	1:C:526:ASP:OD1	2.11	0.51
1:C:342:VAL:HG11	1:D:340:ALA:HB1	1.93	0.51
1:D:685:THR:O	1:D:689:ILE:HG13	2.11	0.51
1:B:285:TYR:CE1	1:B:402:THR:HG22	2.46	0.51
1:B:345:TYR:CD1	1:B:350:LEU:HD23	2.47	0.50
1:A:264:ASP:O	1:A:268:GLN:HG3	2.11	0.50
1:D:354:VAL:O	1:D:354:VAL:HG23	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:667:PRO:O	1:D:671:ARG:HG3	2.11	0.50
1:A:407:TRP:CH2	4:A:1003:HFG:C5'	2.88	0.50
1:A:636:MET:O	1:A:639:THR:HB	2.12	0.50
1:C:267:THR:O	1:C:271:VAL:HG23	2.12	0.50
1:C:406:LEU:HD21	1:D:315:TYR:CE2	2.47	0.50
1:B:409:GLU:HG3	1:B:409:GLU:O	2.12	0.49
1:B:606:VAL:O	1:B:608:SER:O	2.30	0.49
1:C:283:GLY:HA3	1:C:403:ARG:HG3	1.94	0.49
1:A:293:ARG:HH11	1:A:293:ARG:CG	2.25	0.49
1:B:573:VAL:HG22	1:B:574:LYS:H	1.78	0.49
1:D:321:SER:OG	1:D:324:LYS:HG3	2.12	0.49
1:B:669:LEU:HB3	1:B:748:ARG:HH12	1.76	0.49
1:C:417:GLU:HB2	1:C:505:HIS:ND1	2.27	0.49
1:C:475:GLN:NE2	1:C:478:THR:CG2	2.76	0.49
1:C:324:LYS:HD2	1:D:280:ASP:OD2	2.12	0.49
1:D:401:ARG:NH2	2:D:1001:ANP:O3G	2.46	0.49
1:C:365:TYR:N	1:C:365:TYR:CD1	2.80	0.49
1:C:716:LYS:HG2	1:C:717:PRO:O	2.13	0.49
1:D:606:VAL:O	1:D:608:SER:O	2.31	0.49
1:C:454:PHE:CZ	4:C:1003:HFG:H1	2.48	0.48
1:D:364:MET:SD	1:D:411:HIS:CD2	3.06	0.48
1:A:293:ARG:HD3	1:A:537:ALA:CA	2.43	0.48
1:C:587:PHE:O	1:C:591:GLU:HG3	2.13	0.48
1:D:656:ILE:O	1:D:656:ILE:HD12	2.14	0.48
1:B:375:HIS:NE2	1:B:376:ARG:HG3	2.29	0.48
1:C:373:ARG:O	1:C:494:PHE:HA	2.14	0.48
1:D:341:TRP:CE3	1:D:353:LYS:HD2	2.49	0.48
1:D:586:LYS:O	1:D:587:PHE:C	2.51	0.48
1:A:432:ARG:HD3	1:A:445:LYS:HE3	1.94	0.48
1:C:606:VAL:O	1:C:608:SER:O	2.31	0.48
1:D:410:GLY:O	1:D:508:SER:HA	2.12	0.48
1:D:721:PRO:O	1:D:724:GLN:HG2	2.13	0.48
1:D:300:LYS:HE3	1:D:304:ASP:OD2	2.13	0.48
1:C:414:HIS:ND1	1:C:419:GLU:HG2	2.26	0.47
2:C:1001:ANP:O2B	4:C:1003:HFG:H17	2.14	0.47
1:D:493:GLU:HA	1:D:502:GLN:O	2.13	0.47
1:A:265:TRP:O	1:A:268:GLN:HB2	2.14	0.47
1:A:716:LYS:HB2	1:A:717:PRO:HD2	1.96	0.47
1:C:586:LYS:O	1:C:587:PHE:C	2.53	0.47
1:A:336:LYS:N	1:A:337:PRO:HD2	2.29	0.47
1:C:587:PHE:CD1	1:C:598:ARG:HD2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:479:SER:HA	1:D:509:TRP:HB3	1.95	0.47
1:A:410:GLY:O	1:A:508:SER:HA	2.15	0.47
1:B:542:ILE:HB	1:B:575:VAL:HG22	1.95	0.47
1:A:266:TYR:HA	1:A:269:VAL:HG22	1.96	0.47
1:D:300:LYS:O	1:D:300:LYS:HG3	2.14	0.47
1:A:255:THR:H	1:A:268:GLN:HE22	1.62	0.46
1:A:615:ARG:NH2	1:A:617:ASP:OD2	2.48	0.46
1:C:461:THR:OG1	1:C:479:SER:O	2.33	0.46
1:C:533:PRO:HD2	1:C:641:GLN:OE1	2.16	0.46
1:D:424:VAL:HG12	1:D:425:LEU:HD23	1.97	0.46
1:D:513:THR:O	1:D:514:ARG:C	2.53	0.46
1:B:555:ILE:HG13	1:B:606:VAL:HG21	1.97	0.46
1:A:487:SER:OG	1:A:504:VAL:CG2	2.63	0.46
1:B:633:ILE:HB	1:B:634:PRO:HD3	1.97	0.46
1:D:487:SER:O	1:D:491:SER:N	2.48	0.46
1:D:599:LEU:HG	1:D:613:VAL:HG22	1.97	0.46
1:D:279:TYR:CB	1:D:284:CYS:O	2.59	0.46
1:D:725:PRO:HG2	1:D:743:TRP:CG	2.50	0.46
1:B:322:GLN:NE2	1:B:326:GLU:OE2	2.47	0.46
1:A:633:ILE:HB	1:A:634:PRO:HD3	1.97	0.46
1:C:479:SER:HA	1:C:509:TRP:HB3	1.97	0.46
1:C:383:ASN:OD1	1:C:408:GLN:HG2	2.16	0.46
1:B:365:TYR:CD1	1:B:365:TYR:N	2.82	0.46
1:D:365:TYR:CD1	1:D:365:TYR:N	2.83	0.46
1:C:336:LYS:N	1:C:337:PRO:CD	2.78	0.46
1:D:518:VAL:O	1:D:522:THR:OG1	2.28	0.46
1:B:293:ARG:CG	1:B:293:ARG:NH1	2.73	0.45
1:B:660:THR:HG22	1:B:660:THR:O	2.15	0.45
1:D:598:ARG:C	1:D:599:LEU:HD12	2.37	0.45
1:B:316:PHE:HB2	1:B:384:GLN:HE22	1.81	0.45
1:D:261:ASN:OD1	1:D:264:ASP:HB2	2.16	0.45
1:D:527:LYS:O	1:D:592:LEU:HD23	2.16	0.45
1:D:543:ILE:HD12	1:D:598:ARG:CD	2.46	0.45
1:A:320:VAL:CG1	1:A:355:ALA:HB3	2.47	0.45
1:C:289:PRO:HG3	1:D:382:LEU:HD22	1.99	0.45
1:D:434:TRP:C	1:D:434:TRP:CD1	2.89	0.45
1:A:682:ASP:C	1:A:682:ASP:OD1	2.55	0.45
1:D:431:TYR:CE1	1:D:511:CYS:HB2	2.51	0.45
1:A:336:LYS:HE2	1:B:347:ASP:OD1	2.17	0.45
1:A:542:ILE:HB	1:A:575:VAL:HG22	1.98	0.45
1:C:725:PRO:HA	1:C:726:PRO:HD3	1.85	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:356:ILE:HD11	1:D:356:ILE:HD11	1.99	0.45
1:A:554:GLU:O	1:A:558:LYS:HG2	2.17	0.45
1:A:316:PHE:HB2	1:A:384:GLN:HE22	1.82	0.44
1:D:392:GLU:O	1:D:403:ARG:NH2	2.50	0.44
1:C:462:THR:OG1	1:C:475:GLN:HG3	2.17	0.44
1:D:385:TRP:CE3	1:D:408:GLN:HB2	2.52	0.44
1:A:541:VAL:HG12	1:A:595:VAL:HG11	2.00	0.44
1:C:269:VAL:O	1:C:273:GLY:N	2.51	0.44
1:B:275:LEU:HD22	1:B:291:ALA:HB2	2.00	0.44
1:D:281:ILE:O	1:D:282:SER:C	2.55	0.44
1:D:345:TYR:CD2	1:D:345:TYR:O	2.70	0.44
1:A:661:THR:HG22	1:A:664:GLU:OE1	2.17	0.44
1:C:717:PRO:HB3	1:C:746:PHE:CE2	2.52	0.44
1:B:293:ARG:HD3	1:B:537:ALA:HA	1.99	0.44
1:C:434:TRP:C	1:C:434:TRP:CD1	2.91	0.44
1:C:608:SER:HB2	1:C:610:GLN:HG2	1.99	0.44
1:D:428:LEU:HD23	1:D:445:LYS:HG2	1.99	0.44
1:A:666:MET:N	1:A:667:PRO:HD2	2.32	0.44
1:C:269:VAL:HG12	1:C:528:GLY:HA2	2.00	0.44
1:C:527:LYS:HD2	1:C:591:GLU:HB3	2.00	0.44
1:D:453:LYS:O	1:D:454:PHE:C	2.55	0.44
1:C:685:THR:O	1:C:689:ILE:HG13	2.17	0.43
1:D:345:TYR:CD1	1:D:350:LEU:HD23	2.53	0.43
1:A:472:ARG:NH1	1:A:749:SER:O	2.49	0.43
4:C:1003:HFG:N1'	4:C:1003:HFG:O4'	2.51	0.43
1:D:285:TYR:CE1	1:D:402:THR:HG22	2.50	0.43
1:D:666:MET:N	1:D:667:PRO:CD	2.81	0.43
1:C:543:ILE:HD12	1:C:598:ARG:HD2	2.00	0.43
1:C:527:LYS:O	1:C:592:LEU:HD23	2.18	0.43
4:B:1003:HFG:O4'	4:B:1003:HFG:N1'	2.51	0.43
1:D:314:CYS:O	1:D:384:GLN:HB3	2.18	0.43
1:A:277:GLU:HB3	1:A:286:ILE:HB	2.01	0.43
1:A:448:LYS:HE2	1:A:475:GLN:OE1	2.18	0.43
1:A:671:ARG:O	1:A:672:ARG:HB2	2.18	0.43
1:B:450:GLU:O	1:B:453:LYS:HG2	2.18	0.43
1:D:606:VAL:C	1:D:608:SER:O	2.57	0.43
1:D:725:PRO:HG2	1:D:743:TRP:NE1	2.33	0.43
1:C:649:LYS:O	1:C:652:PHE:HB3	2.19	0.43
1:B:318:MET:HE2	1:B:318:MET:HB2	1.84	0.43
1:B:605:ASP:HB3	1:B:610:GLN:O	2.18	0.43
1:C:382:LEU:HD22	1:D:289:PRO:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:374:SER:CB	1:D:376[A]:ARG:HG3	2.48	0.43
1:D:374:SER:OG	1:D:376[A]:ARG:HG3	2.18	0.43
1:D:540:ALA:HB3	1:D:573:VAL:HG12	2.01	0.43
1:D:408:GLN:HE22	1:D:510:GLY:HA2	1.82	0.42
1:D:374:SER:HB3	1:D:376[A]:ARG:HG3	2.01	0.42
1:A:547:PHE:CE2	1:A:603:PRO:HG2	2.54	0.42
1:C:429:GLU:OE1	1:C:429:GLU:HA	2.19	0.42
1:D:366:PRO:HD3	1:D:490:PHE:CD2	2.55	0.42
1:D:462:THR:OG1	1:D:475:GLN:HG3	2.19	0.42
1:C:543:ILE:HD12	1:C:598:ARG:CD	2.50	0.42
1:D:407:TRP:CH2	4:D:1003:HFG:H8	2.54	0.42
1:A:293:ARG:NH1	1:A:293:ARG:CG	2.82	0.42
1:A:315:TYR:CE2	1:B:406:LEU:HD21	2.55	0.42
1:D:429:GLU:OE1	1:D:429:GLU:CA	2.56	0.42
1:A:341:TRP:CZ3	1:A:355:ALA:HB2	2.55	0.42
1:B:541:VAL:HG12	1:B:595:VAL:HG11	2.02	0.42
1:D:487:SER:O	1:D:491:SER:HA	2.20	0.42
1:C:407:TRP:CZ3	4:C:1003:HFG:H8	2.55	0.42
1:B:319:PHE:CE1	1:B:342:VAL:HG21	2.55	0.42
1:C:316:PHE:CE2	1:D:286:ILE:HD13	2.55	0.42
1:C:493:GLU:HA	1:C:502:GLN:O	2.18	0.42
1:D:661:THR:O	1:D:664:GLU:OE1	2.38	0.42
1:A:293:ARG:HH11	1:A:293:ARG:HG2	1.85	0.42
1:D:358:PRO:HA	1:D:388:VAL:HG13	2.01	0.42
1:B:313:ASN:HA	1:B:383:ASN:O	2.20	0.42
1:A:288:ARG:HB3	1:A:289:PRO:HD2	2.02	0.41
1:A:659:ILE:CD1	1:A:665:VAL:HG22	2.50	0.41
1:D:269:VAL:HG23	1:D:270:ILE:HG23	2.02	0.41
1:C:532:PRO:HA	1:C:533:PRO:HD3	1.86	0.41
1:D:345:TYR:O	1:D:345:TYR:CG	2.71	0.41
1:A:444:ILE:HG22	1:A:445:LYS:O	2.21	0.41
1:B:666:MET:N	1:B:667:PRO:CD	2.83	0.41
1:B:718:LEU:O	1:B:719:CYS:HB3	2.20	0.41
1:C:680:CYS:O	1:C:681:GLU:CB	2.68	0.41
1:A:659:ILE:CG1	1:A:677:ALA:HB2	2.51	0.41
1:D:448:LYS:HE2	1:D:475:GLN:HE21	1.84	0.41
1:A:285:TYR:HE1	1:A:402:THR:HG22	1.85	0.41
1:A:314:CYS:O	1:A:384:GLN:HB3	2.21	0.41
1:D:356:ILE:O	1:D:358:PRO:HD3	2.20	0.41
1:A:727:MET:O	1:A:728:PRO:C	2.58	0.41
1:B:423:LEU:HD23	1:B:427:ILE:HG12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:410:GLY:O	1:C:508:SER:HA	2.20	0.41
1:D:275:LEU:O	1:D:288:ARG:HG3	2.21	0.41
1:D:373:ARG:HA	1:D:373:ARG:HD3	1.90	0.41
1:D:459:LYS:HG2	1:D:460:THR:N	2.36	0.41
1:D:632:THR:O	1:D:635:ALA:HB3	2.20	0.41
1:C:533:PRO:HA	1:C:536:VAL:HG12	2.01	0.41
1:D:448:LYS:CE	1:D:475:GLN:NE2	2.84	0.41
1:B:501:LYS:O	1:B:502:GLN:HG2	2.21	0.41
1:B:410:GLY:O	1:B:508:SER:HA	2.21	0.41
1:C:257:LYS:HB2	1:C:260:THR:HB	2.01	0.41
1:C:385:TRP:CE3	1:C:408:GLN:HB2	2.55	0.41
1:C:682:ASP:HA	1:C:683:PRO:HD2	1.98	0.41
1:D:448:LYS:HE2	1:D:475:GLN:NE2	2.35	0.41
1:D:717:PRO:HB3	1:D:746:PHE:CE2	2.55	0.41
1:C:493:GLU:HG3	1:C:503:LEU:HD23	2.04	0.41
1:D:442:PRO:HG3	1:D:649:LYS:HA	2.02	0.41
1:C:264:ASP:O	1:C:268:GLN:HG3	2.21	0.40
1:D:373:ARG:O	1:D:494:PHE:HA	2.21	0.40
1:A:531:LEU:HA	1:A:532:PRO:HD3	1.91	0.40
1:A:632:THR:O	1:A:635:ALA:HB3	2.21	0.40
1:D:280:ASP:OD1	1:D:281:ILE:HG13	2.21	0.40
1:D:386:ASN:OD1	1:D:407:TRP:NE1	2.54	0.40
1:D:487:SER:O	1:D:491:SER:CA	2.69	0.40
1:D:690:LYS:O	1:D:694:GLN:CB	2.69	0.40
2:A:1001:ANP:H8	2:A:1001:ANP:H3'	2.04	0.40
1:C:582:THR:HA	1:C:583:PRO:HD3	2.00	0.40
1:A:317:PRO:HG2	1:B:279:TYR:CE2	2.57	0.40
1:B:344:HIS:HD2	1:B:347:ASP:O	2.04	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	463/498 (93%)	431 (93%)	31 (7%)	1 (0%)	47	79
1	B	470/498 (94%)	433 (92%)	37 (8%)	0	100	100
1	C	454/498 (91%)	417 (92%)	36 (8%)	1 (0%)	47	79
1	D	447/498 (90%)	407 (91%)	39 (9%)	1 (0%)	47	79
All	All	1834/1992 (92%)	1688 (92%)	143 (8%)	3 (0%)	47	79

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	PRO
1	C	358	PRO
1	D	358	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	393/429 (92%)	385 (98%)	8 (2%)	55	80
1	B	394/429 (92%)	380 (96%)	14 (4%)	35	67
1	C	369/429 (86%)	358 (97%)	11 (3%)	41	71
1	D	373/429 (87%)	362 (97%)	11 (3%)	42	72
All	All	1529/1716 (89%)	1485 (97%)	44 (3%)	43	72

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	255	THR
1	A	261	ASN
1	A	293	ARG
1	A	335	PHE
1	A	407	TRP
1	A	408	GLN
1	A	509	TRP
1	A	733	CYS

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Mol	Chain	Res	Type
1	B	293	ARG
1	B	300	LYS
1	B	335	PHE
1	B	357	ARG
1	B	361	GLU
1	B	407	TRP
1	B	408	GLN
1	B	489	MET
1	B	509	TRP
1	B	631	SER
1	B	673	HIS
1	B	674	VAL
1	B	733	CYS
1	B	750	TYR
1	C	260	THR
1	C	293	ARG
1	C	321	SER
1	C	357	ARG
1	C	407	TRP
1	C	408	GLN
1	C	509	TRP
1	C	674	VAL
1	C	727	MET
1	C	733	CYS
1	C	750	TYR
1	D	260	THR
1	D	293	ARG
1	D	300	LYS
1	D	321	SER
1	D	376[A]	ARG
1	D	376[B]	ARG
1	D	407	TRP
1	D	408	GLN
1	D	509	TRP
1	D	615	ARG
1	D	734	PHE

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

Mol	Chain	Res	Type
1	A	475	GLN
1	C	610	GLN

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Mol	Chain	Res	Type
1	D	408	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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	ANP	C	1001	3	29,33,33	2.22	6 (20%)	31,52,52	1.23	2 (6%)
4	HFG	C	1003	-	26,26,26	1.06	1 (3%)	22,37,37	0.78	1 (4%)
2	ANP	D	1001	3	29,33,33	1.76	4 (13%)	31,52,52	1.23	4 (12%)
4	HFG	A	1003	-	26,26,26	1.06	1 (3%)	22,37,37	0.79	1 (4%)
2	ANP	A	1001	3	29,33,33	1.60	3 (10%)	31,52,52	0.94	2 (6%)
4	HFG	B	1003	-	26,26,26	1.05	1 (3%)	22,37,37	0.79	1 (4%)
4	HFG	D	1003	-	26,26,26	1.08	1 (3%)	22,37,37	0.72	1 (4%)
2	ANP	B	1001	3	29,33,33	1.61	3 (10%)	31,52,52	1.13	3 (9%)

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	ANP	C	1001	3	-	3/14/38/38	0/3/3/3
4	HFG	C	1003	-	-	0/8/19/19	0/3/3/3
2	ANP	D	1001	3	-	5/14/38/38	0/3/3/3
4	HFG	A	1003	-	-	0/8/19/19	0/3/3/3
2	ANP	A	1001	3	-	5/14/38/38	0/3/3/3
4	HFG	B	1003	-	-	0/8/19/19	0/3/3/3
4	HFG	D	1003	-	-	2/8/19/19	0/3/3/3
2	ANP	B	1001	3	-	6/14/38/38	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1001	ANP	PG-O1G	6.42	1.56	1.46
2	B	1001	ANP	PG-O1G	6.38	1.56	1.46
2	C	1001	ANP	PG-O1G	5.89	1.55	1.46
2	C	1001	ANP	PB-O2B	-5.70	1.41	1.56
2	A	1001	ANP	PB-O1B	4.93	1.54	1.46
2	A	1001	ANP	PB-O2B	-4.79	1.43	1.56
2	C	1001	ANP	PG-O3G	-4.42	1.44	1.56
2	C	1001	ANP	PB-O1B	4.28	1.52	1.46
2	D	1001	ANP	PG-O3G	-3.86	1.46	1.56
2	B	1001	ANP	PG-O3G	-3.85	1.46	1.56
2	A	1001	ANP	O4'-C1'	3.67	1.46	1.41
2	C	1001	ANP	PG-N3B	3.51	1.72	1.63
4	C	1003	HFG	C4-C10	3.17	1.47	1.41
4	A	1003	HFG	C4-C10	3.16	1.47	1.41
4	B	1003	HFG	C4-C10	3.13	1.47	1.41
2	D	1001	ANP	O4'-C1'	3.12	1.45	1.41
2	D	1001	ANP	PG-N3B	2.75	1.70	1.63
4	D	1003	HFG	C4-C10	2.71	1.46	1.41
2	C	1001	ANP	PB-N3B	2.64	1.70	1.63
2	B	1001	ANP	O4'-C1'	2.23	1.44	1.41

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	ANP	O2G-PG-O1G	-3.12	105.60	113.45
2	B	1001	ANP	O2G-PG-O1G	-3.12	105.62	113.45
2	C	1001	ANP	O2G-PG-O1G	-2.87	106.24	113.45

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	ANP	O1G-PG-N3B	-2.87	107.55	111.77
2	D	1001	ANP	C4-C5-N7	-2.86	106.42	109.40
2	A	1001	ANP	O1G-PG-N3B	-2.58	107.97	111.77
4	B	1003	HFG	C10-C4-N3	-2.25	114.62	116.15
2	D	1001	ANP	O1B-PB-N3B	2.24	115.07	111.77
4	A	1003	HFG	C10-C4-N3	-2.23	114.63	116.15
4	C	1003	HFG	C10-C4-N3	-2.23	114.63	116.15
2	B	1001	ANP	C3'-C2'-C1'	2.20	104.29	100.98
2	A	1001	ANP	O1B-PB-N3B	2.07	114.83	111.77
2	B	1001	ANP	O1B-PB-N3B	2.07	114.81	111.77
4	D	1003	HFG	C10-C4-N3	-2.05	114.75	116.15
2	D	1001	ANP	C3'-C2'-C1'	2.04	104.04	100.98

There are no chirality outliers.

All (21) torsion outliers are listed below:

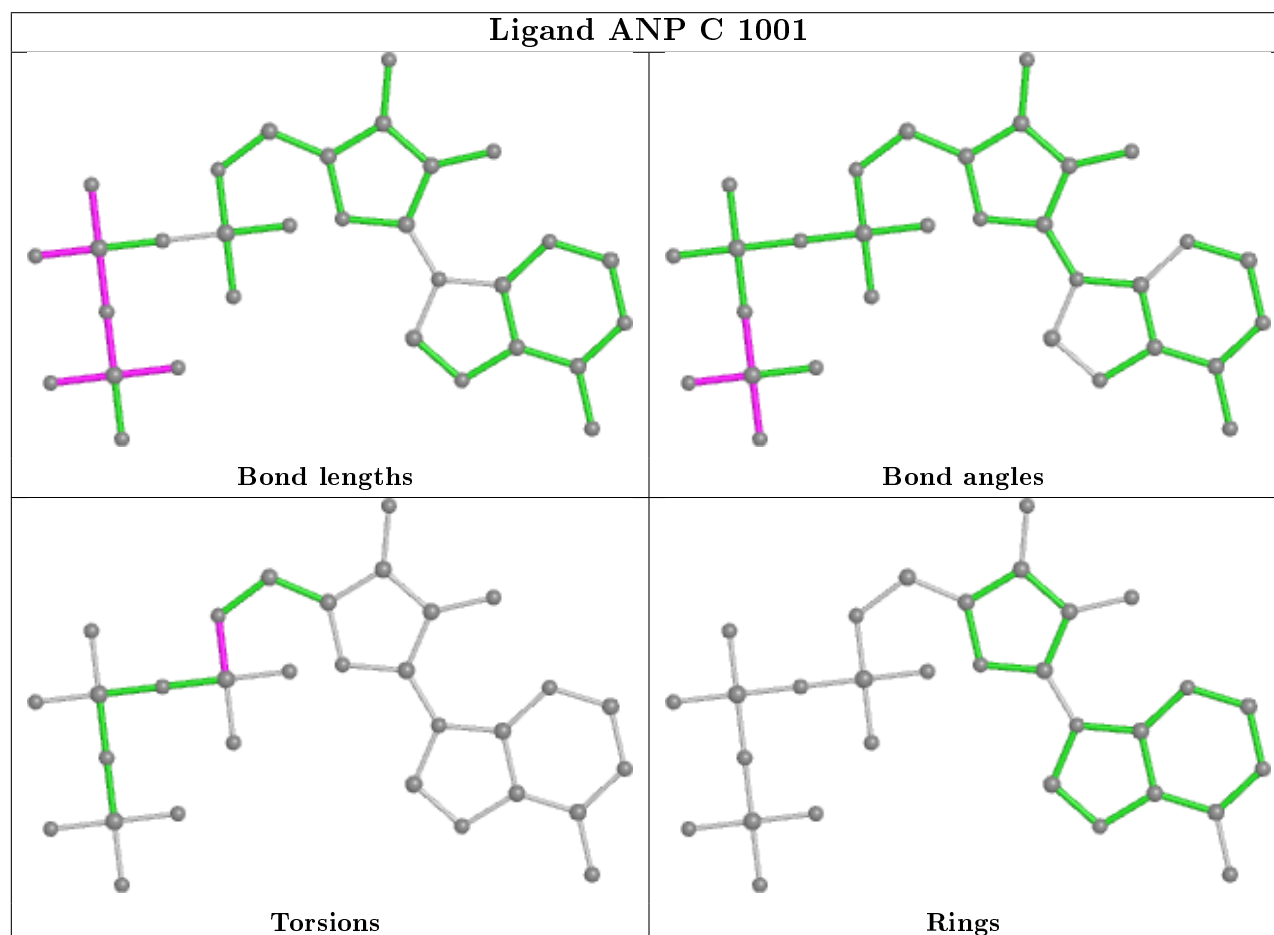
Mol	Chain	Res	Type	Atoms
2	C	1001	ANP	C5'-O5'-PA-O1A
2	C	1001	ANP	C5'-O5'-PA-O2A
2	D	1001	ANP	PB-N3B-PG-O1G
2	D	1001	ANP	PG-N3B-PB-O1B
2	D	1001	ANP	C5'-O5'-PA-O1A
2	D	1001	ANP	C5'-O5'-PA-O2A
2	A	1001	ANP	PB-N3B-PG-O1G
2	A	1001	ANP	C5'-O5'-PA-O1A
2	A	1001	ANP	C5'-O5'-PA-O2A
4	D	1003	HFG	N1'-C2'-C3'-C21
4	D	1003	HFG	C39-C2'-C3'-C21
2	B	1001	ANP	PB-N3B-PG-O1G
2	B	1001	ANP	PG-N3B-PB-O1B
2	B	1001	ANP	PA-O3A-PB-O1B
2	B	1001	ANP	PA-O3A-PB-O2B
2	B	1001	ANP	C5'-O5'-PA-O2A
2	B	1001	ANP	C5'-O5'-PA-O3A
2	A	1001	ANP	C5'-O5'-PA-O3A
2	A	1001	ANP	PB-O3A-PA-O2A
2	C	1001	ANP	C5'-O5'-PA-O3A
2	D	1001	ANP	C5'-O5'-PA-O3A

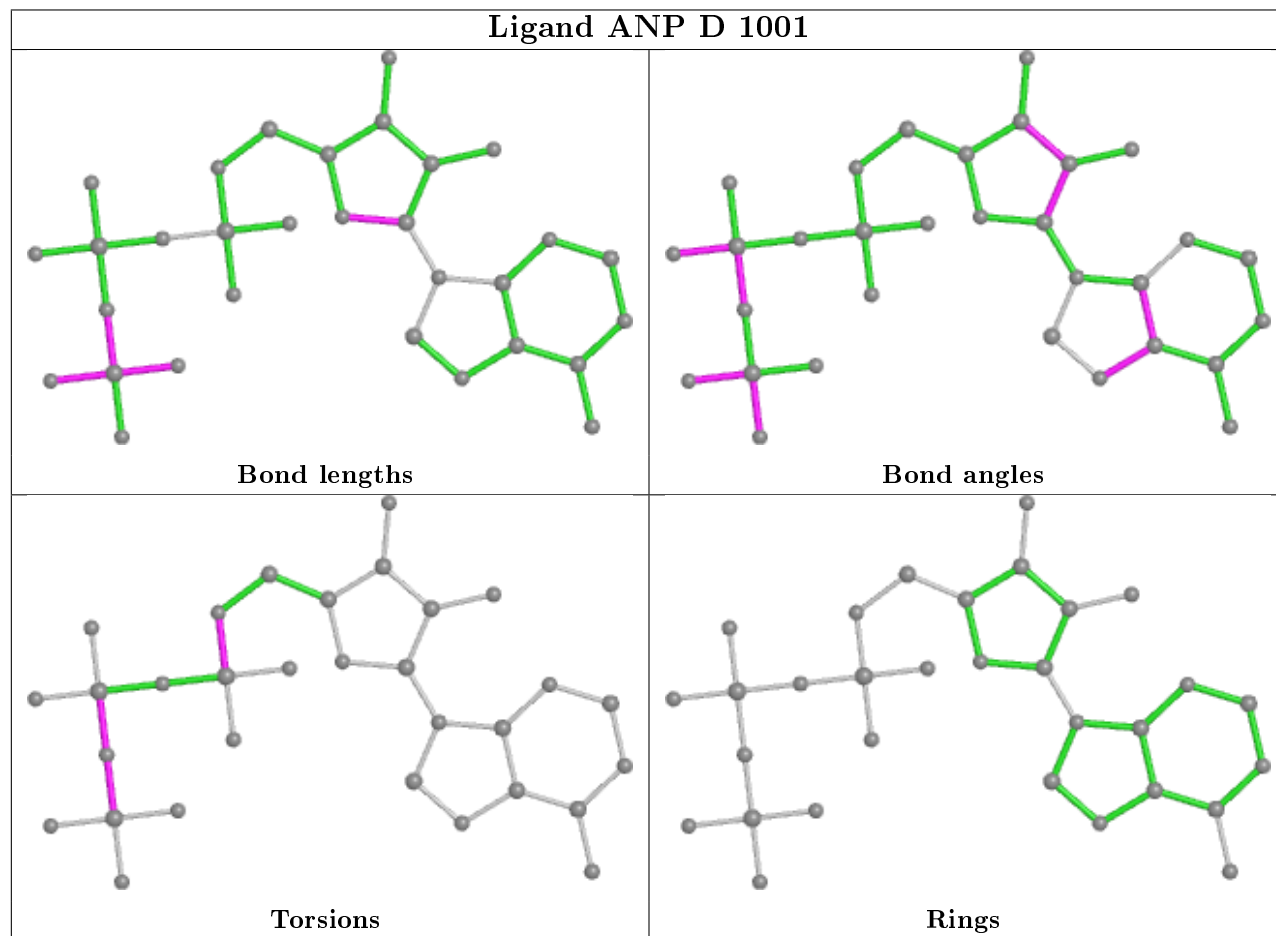
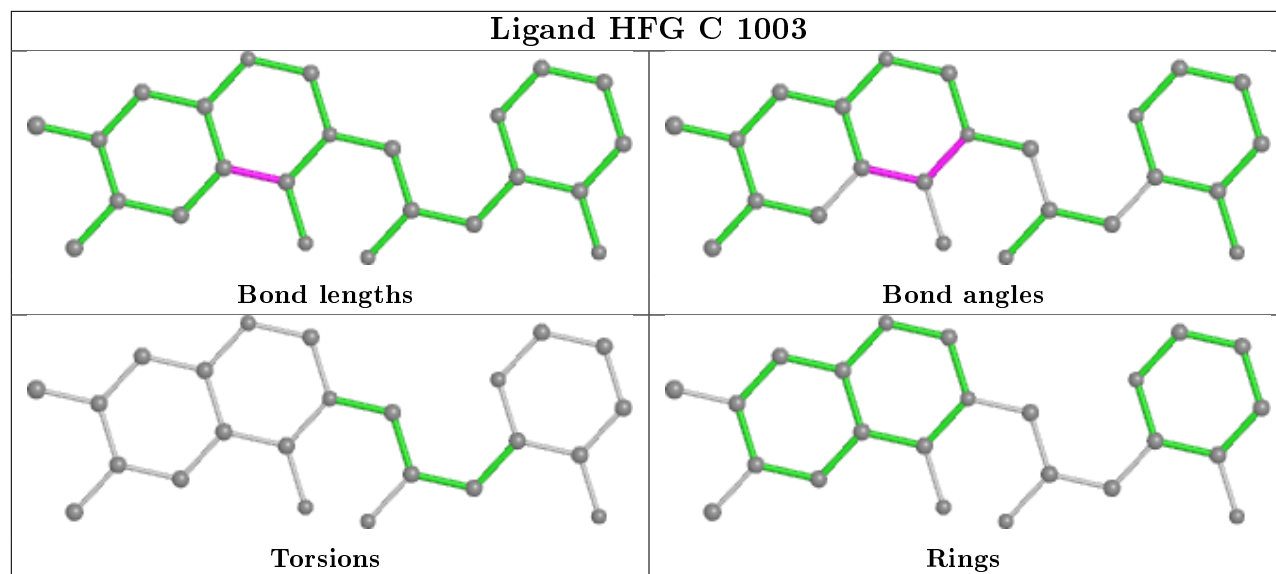
There are no ring outliers.

7 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1001	ANP	2	0
4	C	1003	HFG	6	0
2	D	1001	ANP	4	0
4	A	1003	HFG	6	0
2	A	1001	ANP	2	0
4	B	1003	HFG	7	0
4	D	1003	HFG	4	0

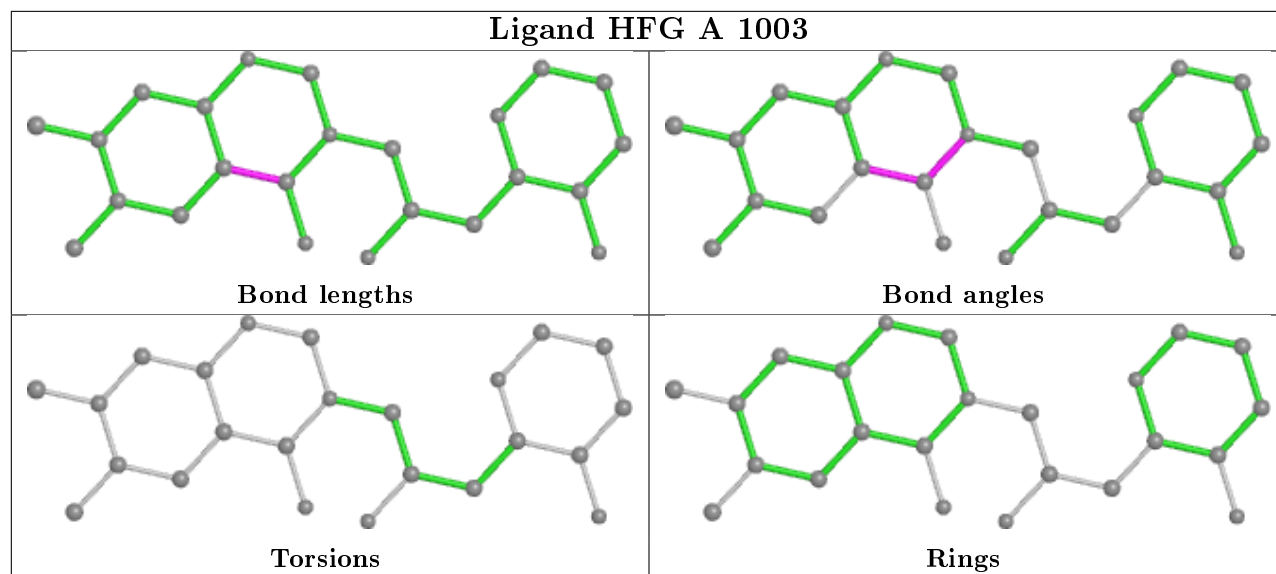
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



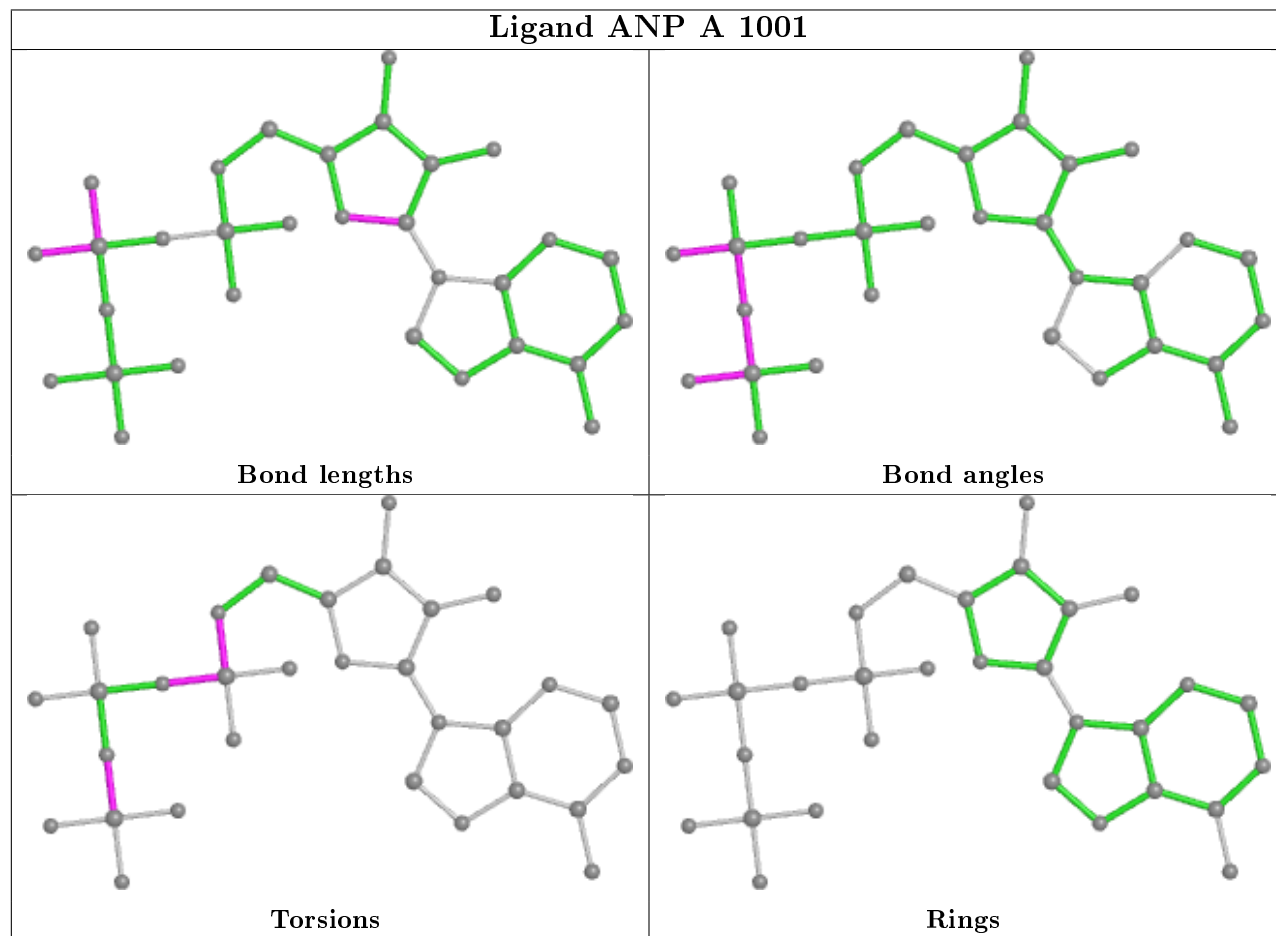


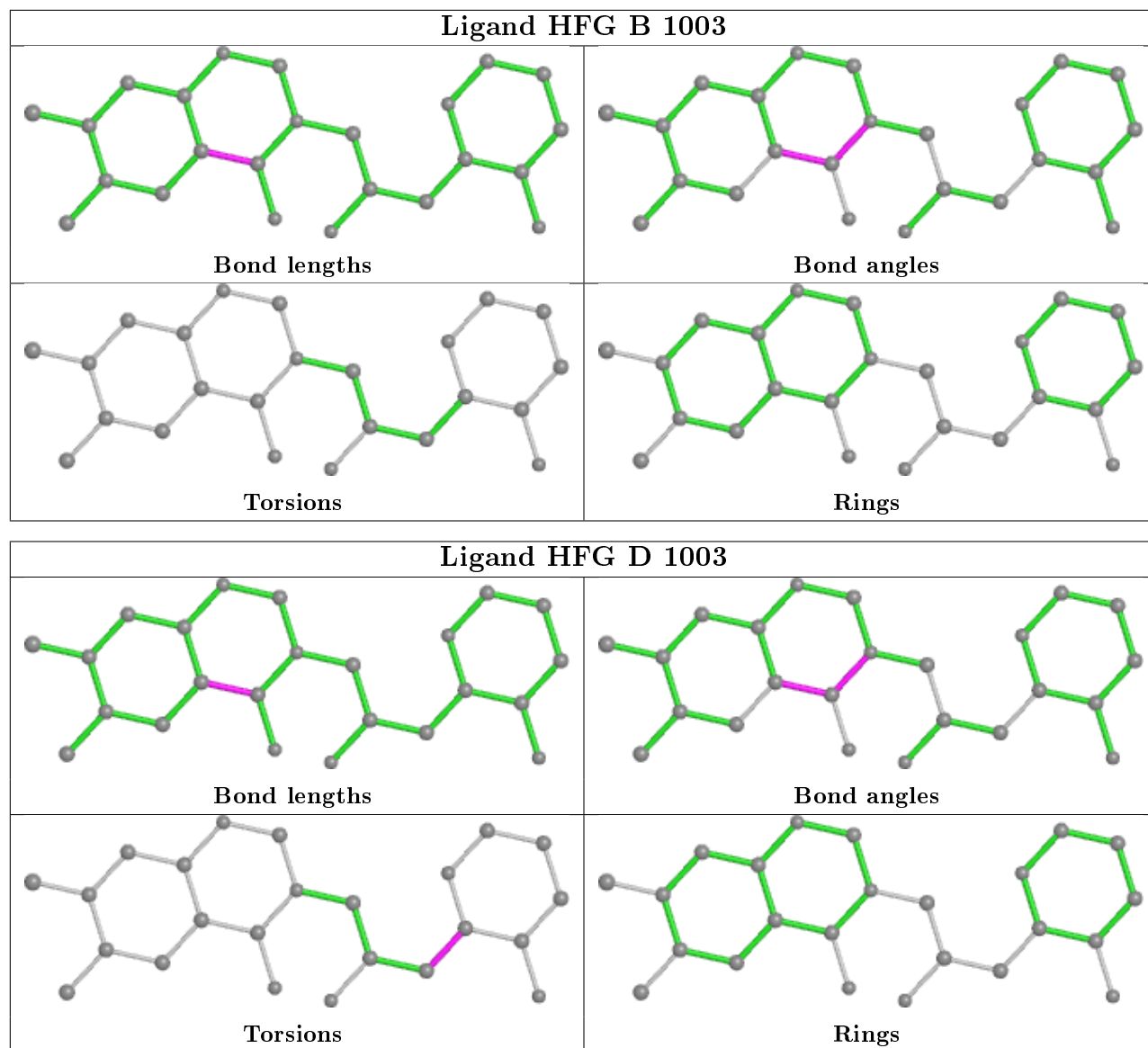


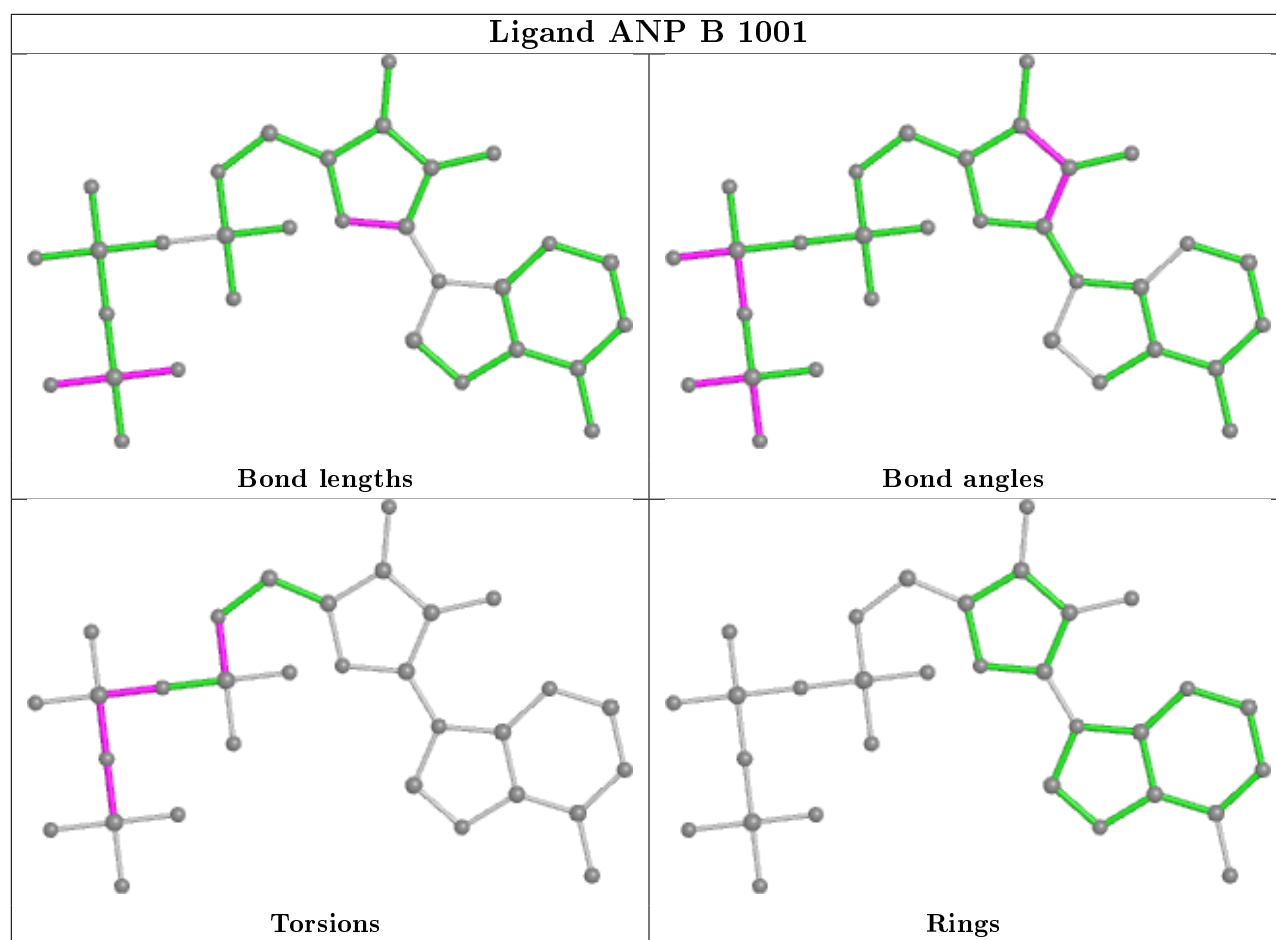
## Ligand HFG A 1003



## Ligand ANP A 1001







## 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	471/498 (94%)	-0.19	0 100 100	10, 26, 45, 59	0
1	B	475/498 (95%)	-0.18	1 (0%) 95 90	11, 26, 48, 71	0
1	C	463/498 (92%)	-0.10	2 (0%) 92 84	12, 34, 69, 80	0
1	D	457/498 (91%)	-0.11	0 100 100	11, 33, 68, 84	0
All	All	1866/1992 (93%)	-0.14	3 (0%) 95 90	10, 28, 63, 84	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	345	TYR	2.4
1	B	335	PHE	2.1
1	C	659	ILE	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](#)

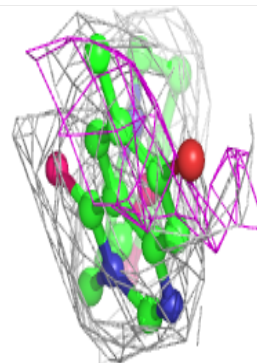
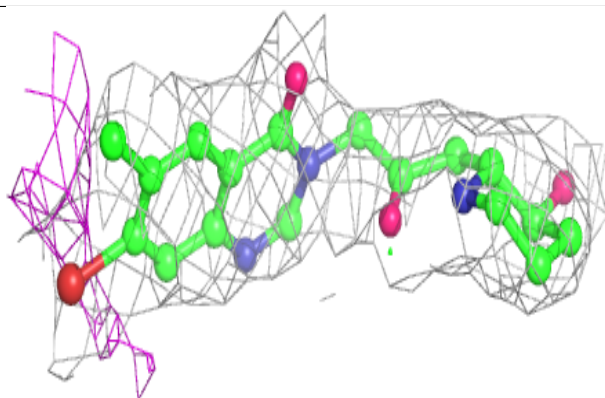
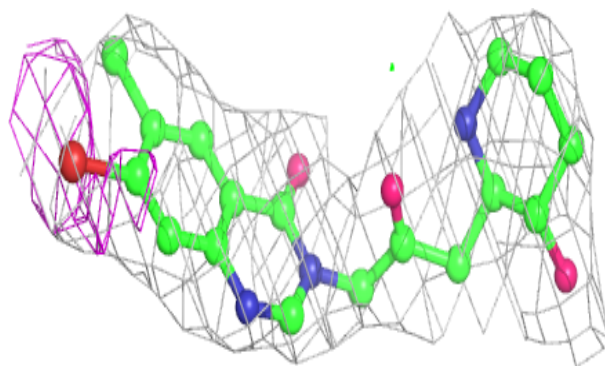
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 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
3	MG	B	1002	1/1	0.86	0.11	33,33,33,33	0
3	MG	C	1002	1/1	0.88	0.12	31,31,31,31	0
4	HFG	D	1003	24/24	0.90	0.23	47,54,70,98	0
4	HFG	B	1003	24/24	0.91	0.24	29,34,48,75	0
4	HFG	C	1003	24/24	0.91	0.20	39,51,63,84	0
2	ANP	C	1001	31/31	0.94	0.17	34,43,49,50	0
2	ANP	D	1001	31/31	0.94	0.17	36,41,44,46	0
4	HFG	A	1003	24/24	0.94	0.20	21,31,40,50	0
2	ANP	A	1001	31/31	0.96	0.19	14,18,23,24	0
2	ANP	B	1001	31/31	0.96	0.17	14,20,27,28	0
3	MG	D	1002	1/1	0.98	0.06	44,44,44,44	0
3	MG	A	1002	1/1	0.98	0.14	18,18,18,18	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.

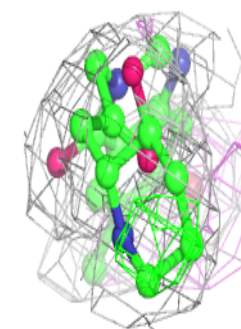
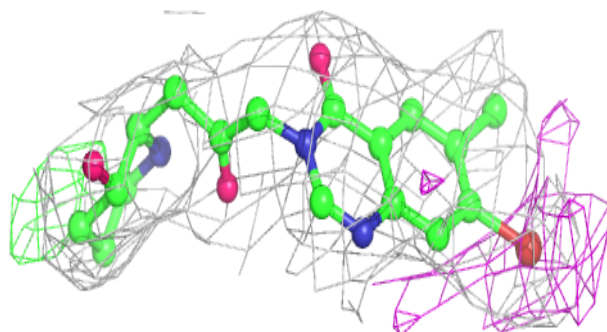
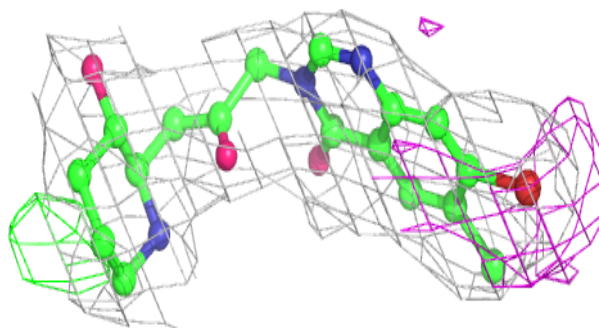
**Electron density around HFG D 1003:**

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)

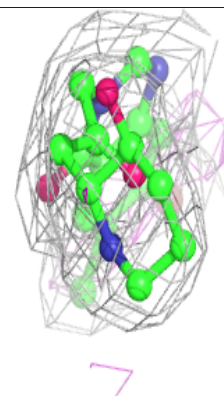
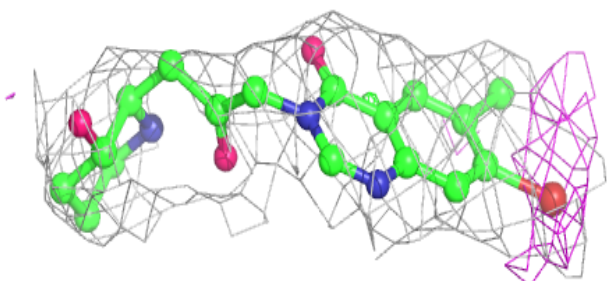
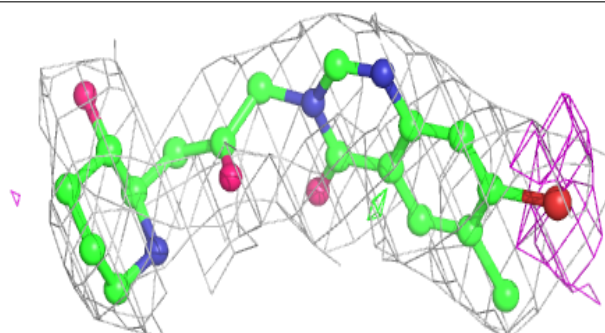


**Electron density around HFG B 1003:**

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

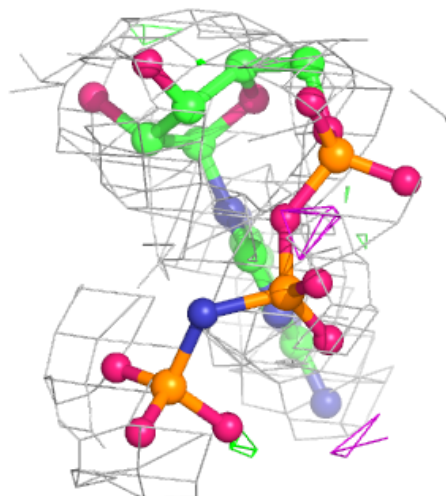
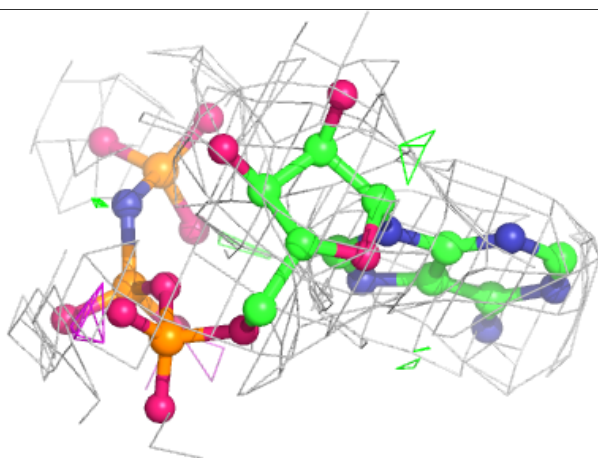
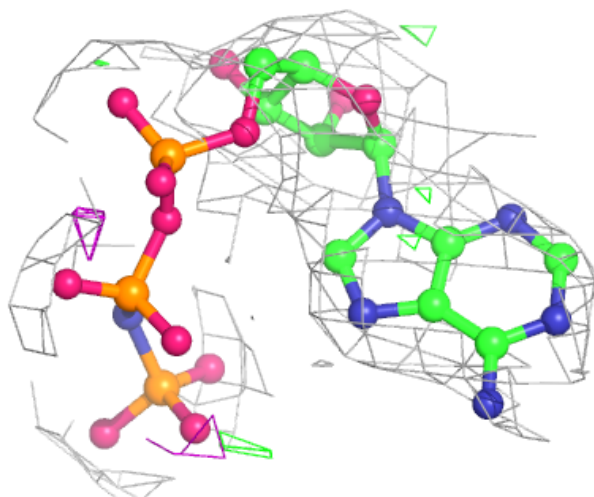
**Electron density around HFG C 1003:**

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



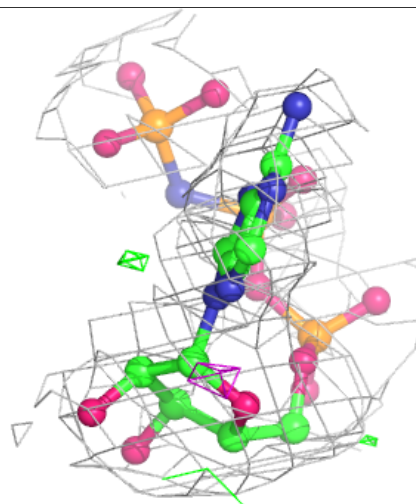
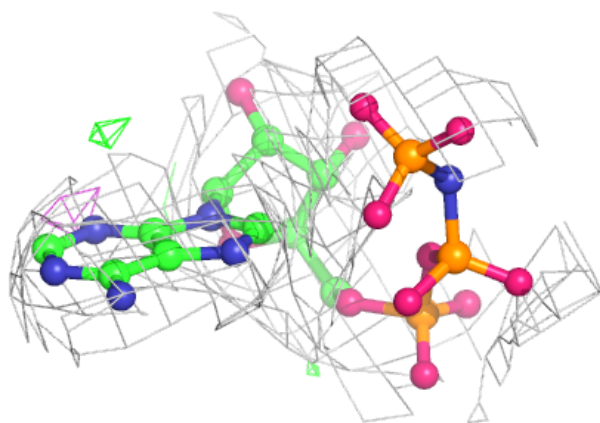
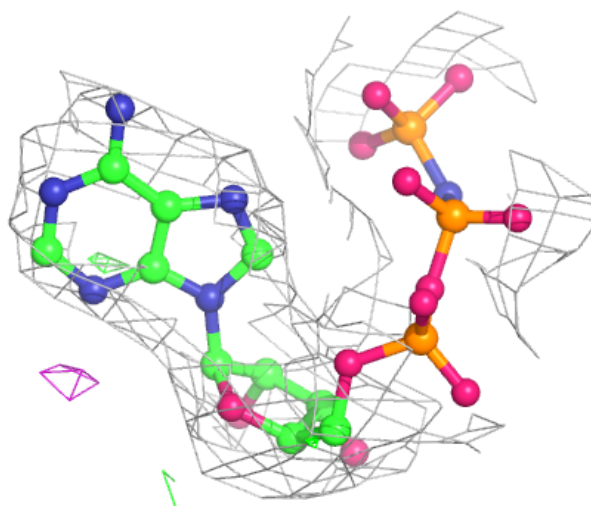
**Electron density around ANP C 1001:**

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



**Electron density around ANP D 1001:**

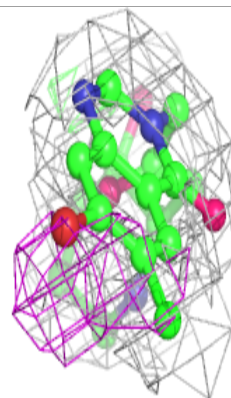
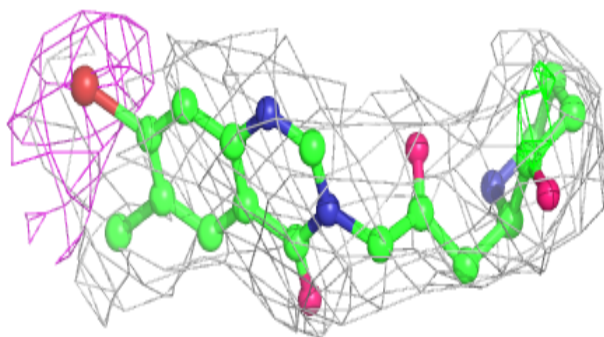
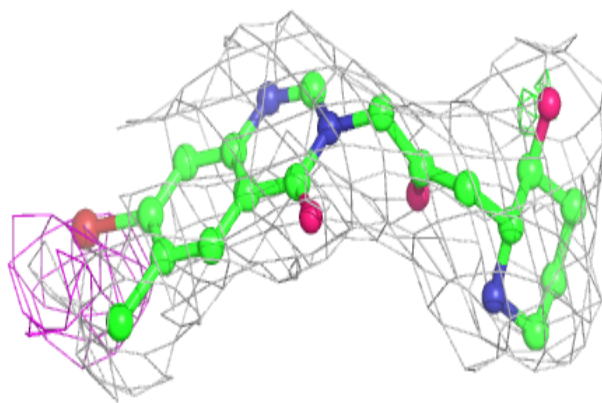
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





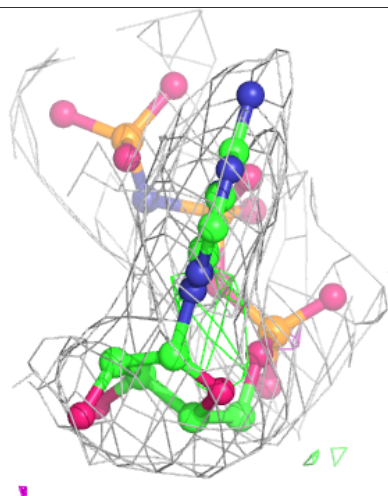
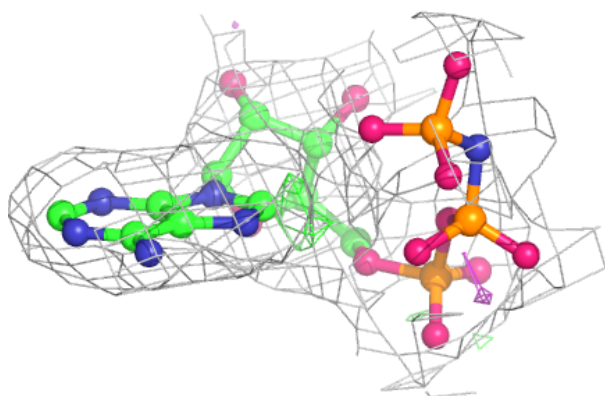
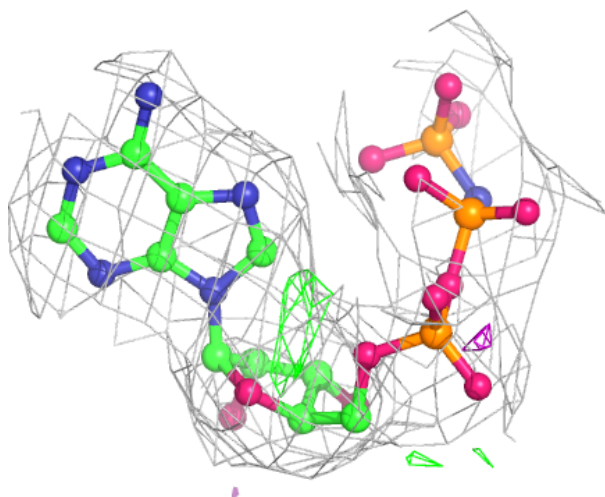
**Electron density around HFG A 1003:**

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



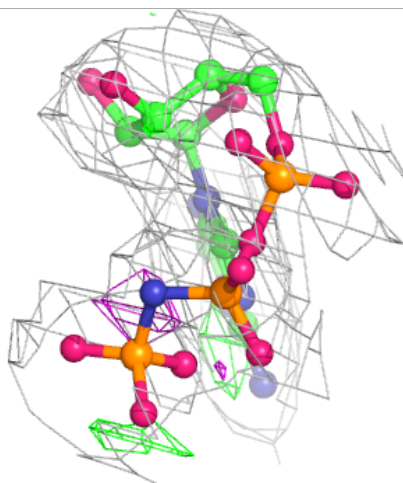
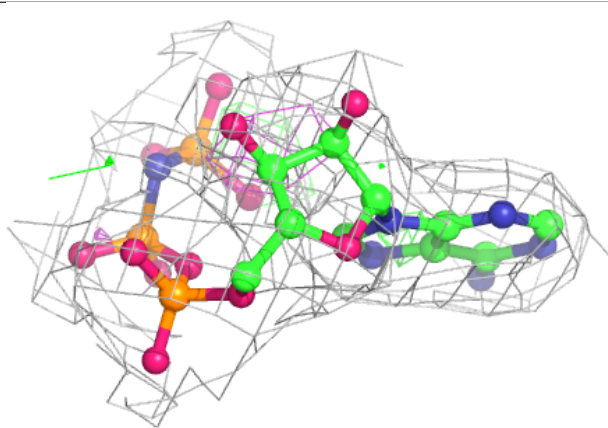
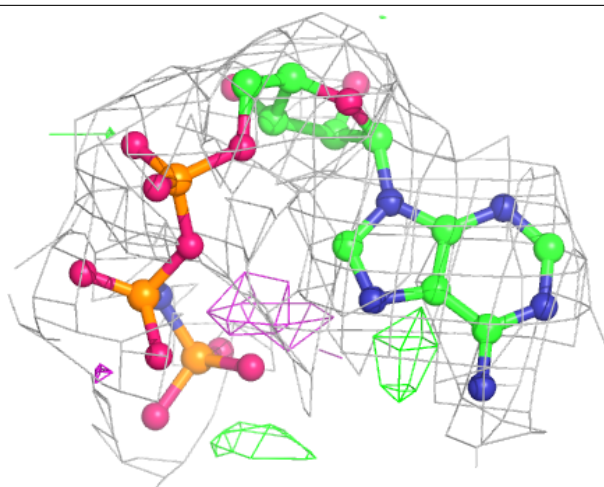
**Electron density around ANP A 1001:**

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



**Electron density around ANP B 1001:**

$2mF_o-DF_c$  (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.