



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:05 AM UTC

PDB ID : 2XND / pdb_00002xnd
Title : Crystal structure of bovine F1-c8 sub-complex of ATP Synthase
Authors : Watt, I.N.; Montgomery, M.G.; Runswick, M.J.; Leslie, A.G.W.; Walker, J.E.
Deposited on : 2010-08-02
Resolution : 3.50 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

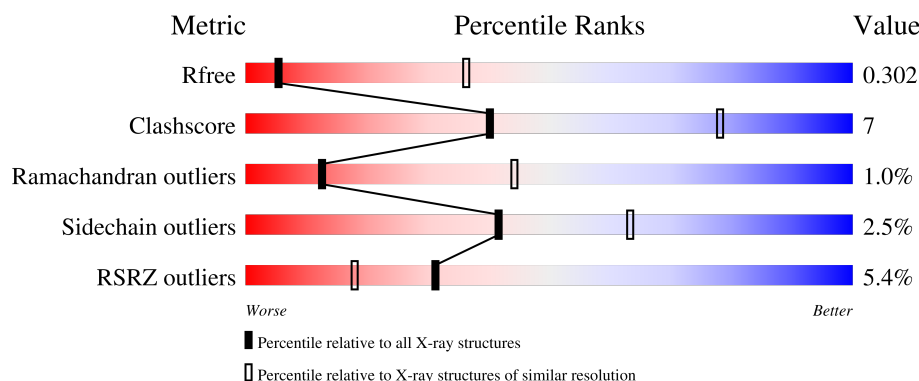
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	492	
1	B	492	
1	C	492	
2	D	467	
2	E	467	

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Mol	Chain	Length	Quality of chain
2	F	467	
3	G	272	
4	H	131	
5	I	47	
6	J	72	
6	K	72	
6	L	72	
6	M	72	
6	N	72	
6	O	72	
6	P	72	
6	Q	72	

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 28196 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 ATP SYNTHASE SUBUNIT ALPHA, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	0	0
			3748	2360	661	715	12			
1	B	492	Total	C	N	O	S	0	0	0
			3748	2360	661	715	12			
1	C	492	Total	C	N	O	S	0	0	0
			3748	2360	661	715	12			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	GLY	SER	cloning artifact	UNP P19483
B	481	GLY	SER	cloning artifact	UNP P19483
C	481	GLY	SER	cloning artifact	UNP P19483

- Molecule 2 is a protein called ATP SYNTHASE SUBUNIT BETA, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	467	Total	C	N	O	S	0	0	0
			3539	2243	601	684	11			
2	E	466	Total	C	N	O	S	0	0	0
			3530	2238	600	681	11			
2	F	466	Total	C	N	O	S	0	0	0
			3530	2238	600	681	11			

- Molecule 3 is a protein called ATP SYNTHASE SUBUNIT GAMMA, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	263	Total	C	N	O	S	0	0	0
			2051	1291	354	398	8			

- Molecule 4 is a protein called ATP SYNTHASE SUBUNIT DELTA, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	131	Total	C	N	O	S	0	0	0
			970	609	164	195	2			

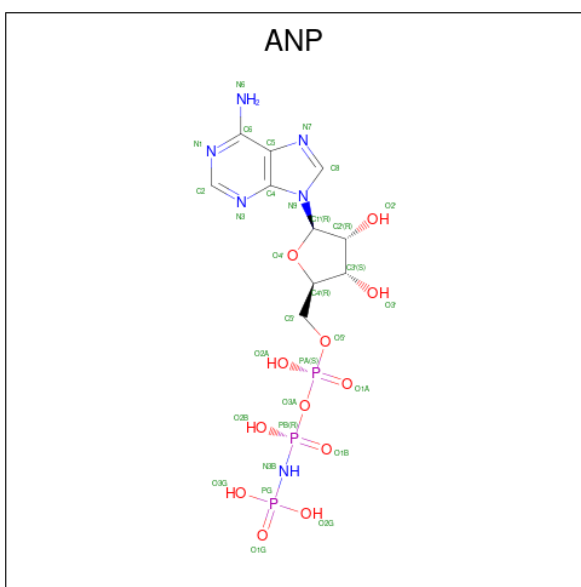
- Molecule 5 is a protein called ATP SYNTHASE SUBUNIT EPSILON, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	47	Total	C	N	O	S	0	0	0
			369	237	66	64	2			

- Molecule 6 is a protein called ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	J	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	K	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	L	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	M	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	N	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	O	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	P	72	Total	C	N	O		0	0	0
			349	205	72	72				
6	Q	72	Total	C	N	O		0	0	0
			349	205	72	72				

- Molecule 7 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).

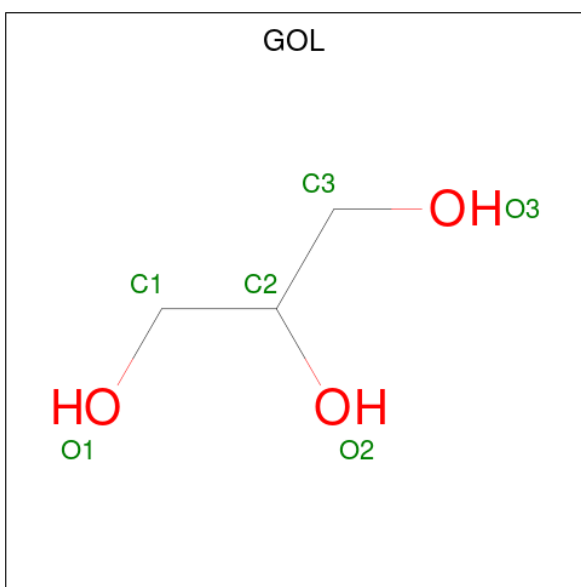


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
7	B	1	Total 31	C 10	N 6	O 12	P 3	0	0
7	C	1	Total 31	C 10	N 6	O 12	P 3	0	0
7	D	1	Total 31	C 10	N 6	O 12	P 3	0	0
7	F	1	Total 31	C 10	N 6	O 12	P 3	0	0

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

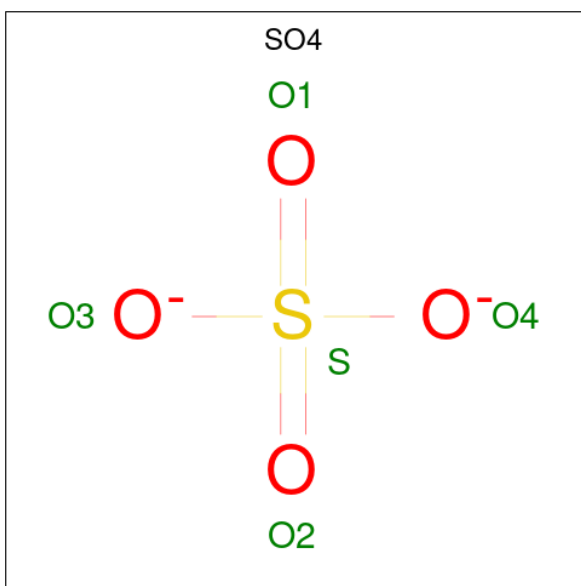
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Mg 1 1	0	0
8	B	1	Total Mg 1 1	0	0
8	C	1	Total Mg 1 1	0	0
8	D	1	Total Mg 1 1	0	0
8	F	1	Total Mg 1 1	0	0

- Molecule 9 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

- Molecule 10 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

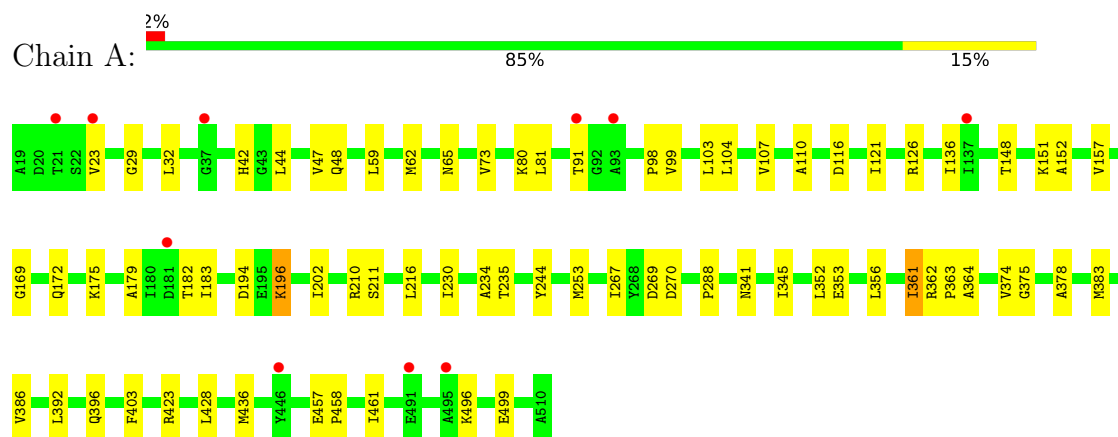


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	E	1	Total	O	S	0	0
			5	4	1		

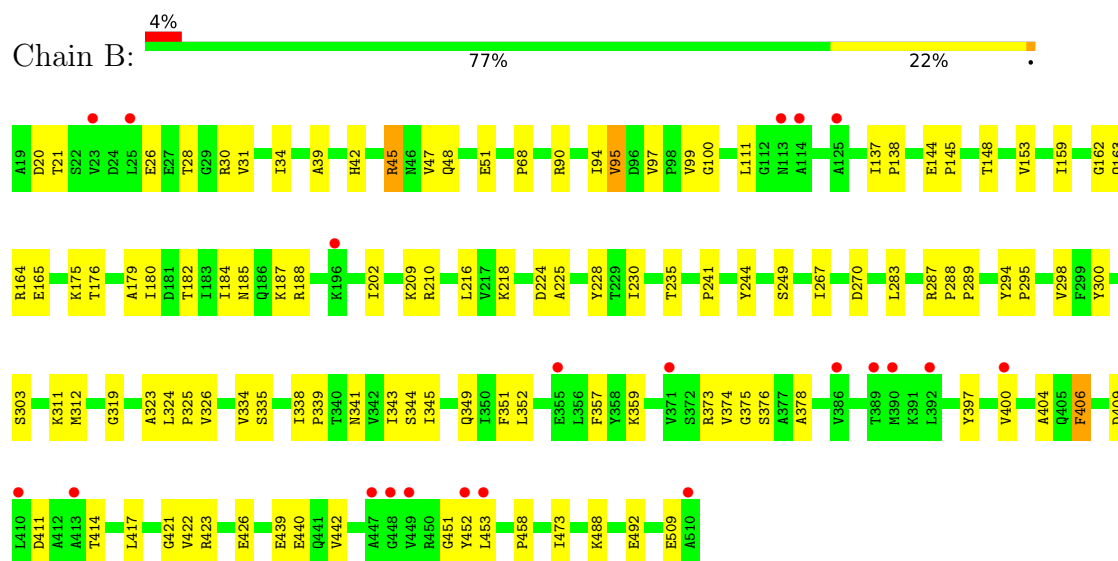
3 Residue-property plots [i](#)

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.

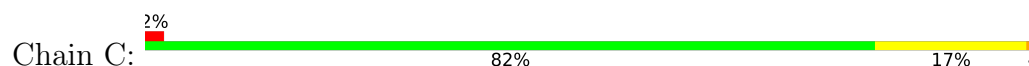
• Molecule 1: ATP SYNTHASE SUBUNIT ALPHA, MITOCHONDRIAL

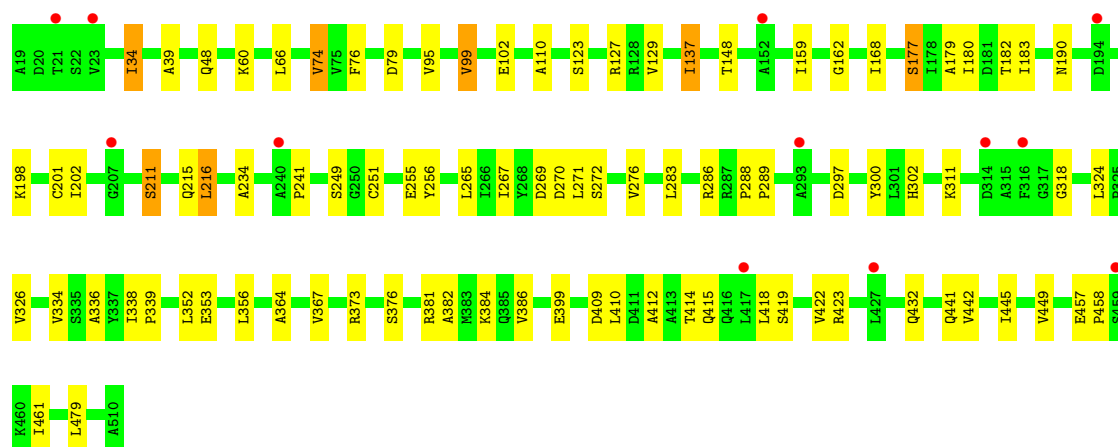


• Molecule 1: ATP SYNTHASE SUBUNIT ALPHA, MITOCHONDRIAL

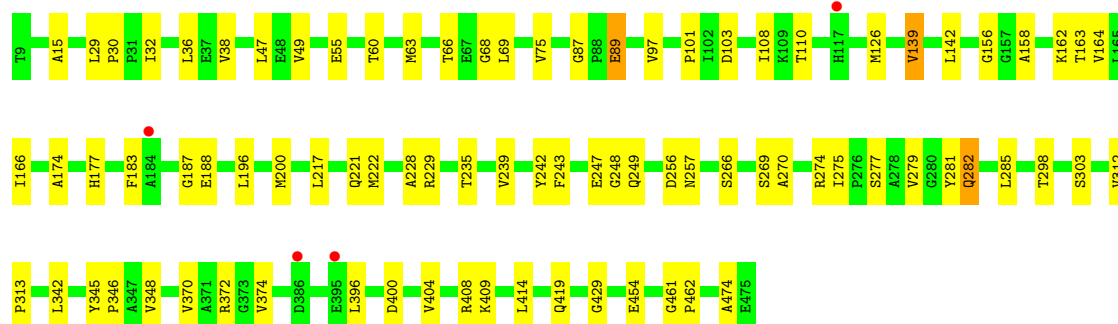
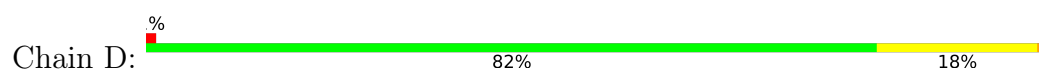


• Molecule 1: ATP SYNTHASE SUBUNIT ALPHA, MITOCHONDRIAL

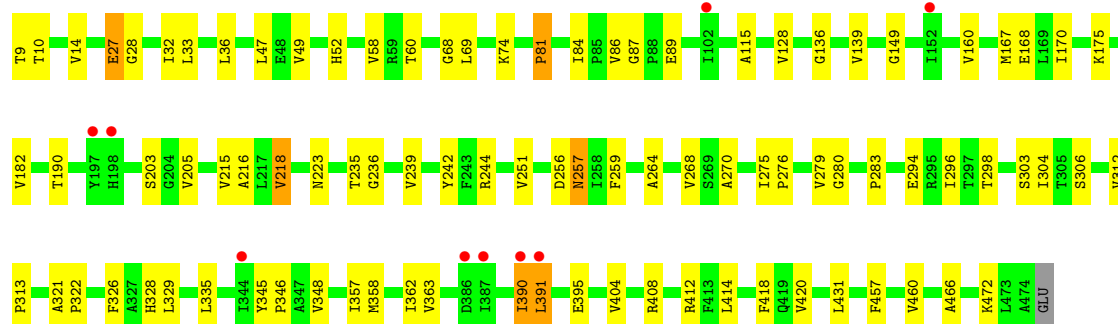
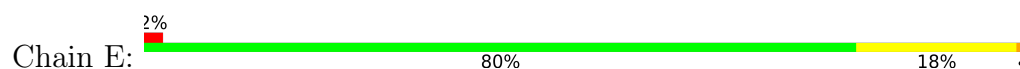




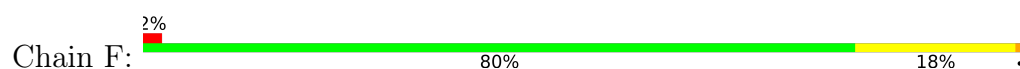
• Molecule 2: ATP SYNTHASE SUBUNIT BETA, MITOCHONDRIAL

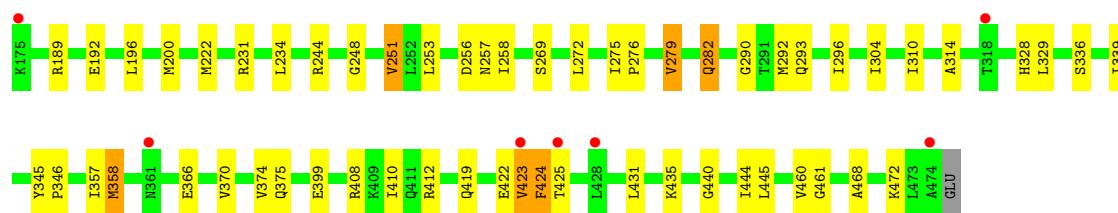


• Molecule 2: ATP SYNTHASE SUBUNIT BETA, MITOCHONDRIAL

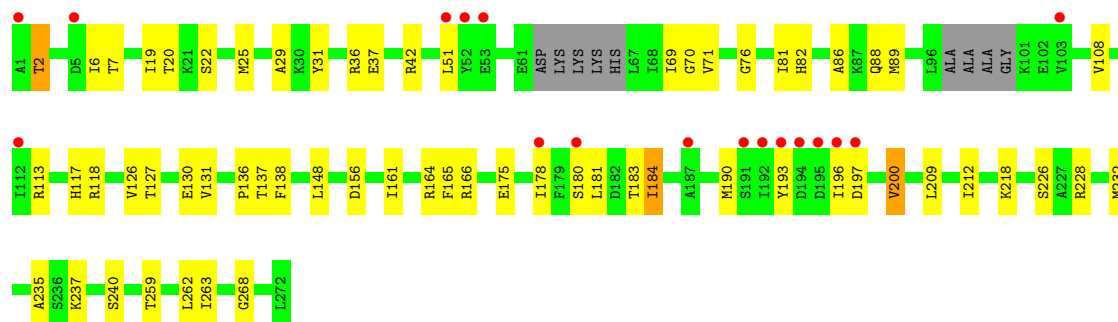
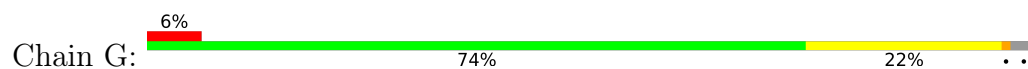


• Molecule 2: ATP SYNTHASE SUBUNIT BETA, MITOCHONDRIAL

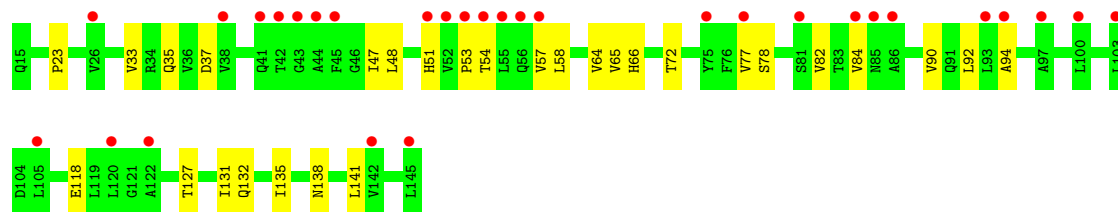
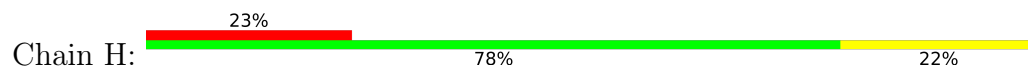




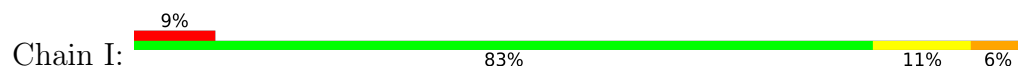
• Molecule 3: ATP SYNTHASE SUBUNIT GAMMA, MITOCHONDRIAL



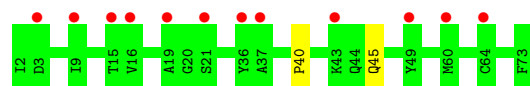
• Molecule 4: ATP SYNTHASE SUBUNIT DELTA, MITOCHONDRIAL



• Molecule 5: ATP SYNTHASE SUBUNIT EPSILON, MITOCHONDRIAL

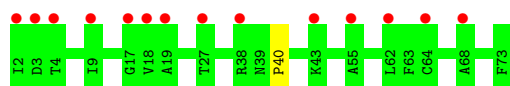


• Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL

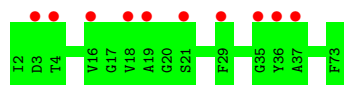


• Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL

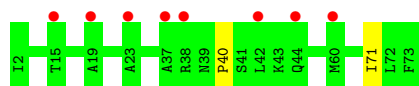




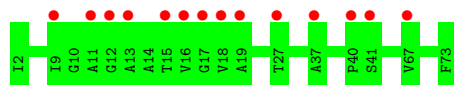
- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



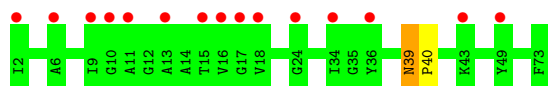
- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



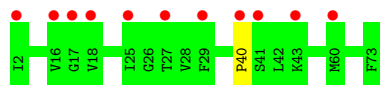
- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



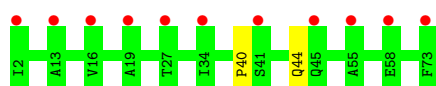
- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



- Molecule 6: ATP SYNTHASE LIPID-BINDING PROTEIN, MITOCHONDRIAL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	155.74Å 157.10Å 247.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	132.60 – 3.50 132.60 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (132.60-3.50) 99.6 (132.60-3.50)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.261 , 0.304 0.262 , 0.302	Depositor DCC
R_{free} test set	3855 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	76.6	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 20.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.088 for k,h,-l	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	28196	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: MG, GOL, ANP, SO4

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	0/3799	0.81	0/5126
1	B	0.45	0/3799	0.79	3/5126 (0.1%)
1	C	0.47	1/3799 (0.0%)	0.81	0/5126
2	D	0.45	0/3596	0.81	0/4879
2	E	0.45	0/3587	0.79	0/4867
2	F	0.46	0/3587	0.82	2/4867 (0.0%)
3	G	0.44	0/2074	0.78	0/2785
4	H	0.49	0/982	0.78	0/1337
5	I	0.49	0/374	0.77	0/501
6	J	0.48	0/348	0.91	0/479
6	K	0.48	0/348	0.89	0/479
6	L	0.49	0/348	0.88	0/479
6	M	0.46	0/348	0.92	0/479
6	N	0.46	0/348	0.89	0/479
6	O	0.47	0/348	0.91	0/479
6	P	0.46	0/348	0.90	0/479
6	Q	0.47	0/348	0.89	0/479
All	All	0.46	1/28381 (0.0%)	0.81	5/38446 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	137	ILE	CA-CB	6.18	1.57	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	87	GLY	CA-C-N	5.39	125.06	119.56
2	F	87	GLY	C-N-CA	5.39	125.06	119.56
1	B	45	ARG	N-CA-C	5.04	118.20	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	287	ARG	CA-C-N	5.03	125.56	120.38
1	B	287	ARG	C-N-CA	5.03	125.56	120.38

There are no chirality outliers.

There are no planarity outliers.

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	3748	0	3844	46	0
1	B	3748	0	3844	65	0
1	C	3748	0	3844	56	0
2	D	3539	0	3592	56	0
2	E	3530	0	3587	55	0
2	F	3530	0	3586	57	0
3	G	2051	0	2115	37	0
4	H	970	0	972	18	0
5	I	369	0	395	6	0
6	J	349	0	189	0	0
6	K	349	0	189	0	0
6	L	349	0	189	0	0
6	M	349	0	189	0	0
6	N	349	0	189	0	0
6	O	349	0	189	1	0
6	P	349	0	189	0	0
6	Q	349	0	189	0	0
7	A	31	0	13	1	0
7	B	31	0	13	1	0
7	C	31	0	13	2	0
7	D	31	0	13	3	0
7	F	31	0	13	4	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
8	F	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	6	0	8	0	0
10	E	5	0	0	0	0
All	All	28196	0	27364	364	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 7.

All (364) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:282:GLN:H	2:F:282:GLN:HE21	1.15	0.93
1:A:99:VAL:HG23	1:A:253:MET:HA	1.57	0.86
2:E:27:GLU:HG2	2:E:28:GLY:H	1.41	0.85
2:F:51:GLN:HE21	2:F:59:ARG:HD2	1.46	0.78
2:E:298:THR:HG23	2:E:303:SER:HB3	1.66	0.77
2:F:89:GLU:HB2	2:F:110:THR:HG22	1.67	0.75
2:E:257:ASN:HD21	2:E:259:PHE:HB3	1.54	0.72
1:A:403:PHE:HE1	3:G:19:ILE:HA	1.55	0.71
2:D:282:GLN:H	2:D:282:GLN:HE21	1.37	0.70
2:D:298:THR:HG23	2:D:303:SER:HA	1.73	0.70
1:C:423:ARG:HD2	1:C:461:ILE:HD11	1.72	0.70
1:A:341:ASN:O	1:A:345:ILE:HG12	1.92	0.69
1:A:44:LEU:O	1:A:47:VAL:HG22	1.94	0.68
1:C:179:ALA:O	1:C:183:ILE:HG12	1.92	0.68
1:C:419:SER:O	1:C:423:ARG:HG2	1.93	0.68
2:E:390:ILE:HG22	2:E:390:ILE:O	1.94	0.67
2:E:136:GLY:HA3	2:E:431:LEU:HD22	1.77	0.66
3:G:148:LEU:HD23	3:G:209:LEU:HD21	1.77	0.66
1:A:169:GLY:H	1:A:175:LYS:HD3	1.61	0.66
1:C:74:VAL:CG1	1:C:241:PRO:HB3	2.26	0.66
1:B:339:PRO:O	1:B:343:ILE:HG12	1.96	0.65
3:G:71:VAL:HA	3:G:108:VAL:HG12	1.78	0.65
3:G:156:ASP:HA	3:G:181:LEU:HD22	1.79	0.64
1:C:99:VAL:HG21	1:C:127:ARG:HB3	1.78	0.64
2:D:139:VAL:HG11	2:D:348:VAL:HB	1.80	0.63
1:C:48:GLN:HB3	2:D:68:GLY:HA2	1.80	0.63
1:B:343:ILE:HB	2:F:158:ALA:HB1	1.81	0.63
1:C:168:ILE:HD11	1:C:339:PRO:HB3	1.79	0.63
1:B:99:VAL:HG12	1:B:100:GLY:N	2.12	0.63
1:B:406:PHE:HB2	1:B:409:ASP:HB2	1.80	0.63
2:E:390:ILE:HG23	3:G:29:ALA:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:290:GLY:HA2	2:F:328:HIS:HE1	1.64	0.63
4:H:54:THR:H	4:H:84:VAL:HG23	1.64	0.62
2:D:188:GLU:O	2:D:221:GLN:HB3	1.99	0.62
2:F:419:GLN:HE21	2:F:431:LEU:HD13	1.65	0.62
1:C:441:GLN:O	1:C:445:ILE:HG12	2.01	0.61
2:E:390:ILE:O	2:E:390:ILE:CG2	2.48	0.61
2:D:29:LEU:HD12	2:D:30:PRO:HD2	1.81	0.61
4:H:47:ILE:HG21	4:H:84:VAL:HG11	1.82	0.61
1:A:211:SER:HB3	2:D:126:MET:HE2	1.82	0.61
3:G:197:ASP:HB2	3:G:200:VAL:HB	1.81	0.61
1:C:99:VAL:HG13	1:C:256:TYR:HB2	1.81	0.60
1:B:99:VAL:HG12	1:B:100:GLY:H	1.66	0.60
2:D:419:GLN:HA	2:D:429:GLY:HA3	1.84	0.60
1:A:210:ARG:HG3	1:A:235:THR:HG21	1.83	0.60
1:B:218:LYS:HG3	2:E:128:VAL:HB	1.83	0.59
1:A:151:LYS:HD2	1:A:436:MET:SD	2.43	0.59
1:C:74:VAL:HG13	1:C:241:PRO:HB3	1.83	0.59
2:D:279:VAL:HG12	2:D:279:VAL:O	2.03	0.59
1:C:129:VAL:HG12	1:C:249:SER:HA	1.83	0.59
2:E:84:ILE:HG21	2:E:235:THR:HG23	1.84	0.58
1:A:179:ALA:HB1	1:A:267:ILE:HD13	1.85	0.58
2:F:147:ALA:HB2	2:F:357:ILE:HG13	1.85	0.58
1:B:283:LEU:HD21	1:B:289:PRO:HB3	1.84	0.58
1:A:59:LEU:HD11	1:A:81:LEU:HD13	1.84	0.58
1:C:34:ILE:HD11	1:C:79:ASP:HB2	1.84	0.58
2:E:408:ARG:O	2:E:412:ARG:HG2	2.04	0.58
2:F:162:LYS:HE3	2:F:257:ASN:HD21	1.69	0.58
2:D:139:VAL:HG13	2:D:414:LEU:HD22	1.86	0.57
2:D:183:PHE:HB3	2:D:217:LEU:HD23	1.86	0.57
2:F:189:ARG:HB2	2:F:192:GLU:HG3	1.85	0.57
2:E:167:MET:HA	2:E:170:ILE:HD12	1.85	0.57
1:B:162:GLY:HA3	1:B:311:LYS:HB2	1.86	0.57
3:G:178:ILE:HG22	3:G:180:SER:HB2	1.87	0.56
3:G:259:THR:O	3:G:263:ILE:HG12	2.04	0.56
1:B:34:ILE:HD13	1:B:39:ALA:HB2	1.87	0.56
1:B:335:SER:O	2:F:314:ALA:HA	2.05	0.56
1:B:439:GLU:HA	1:B:442:VAL:HG22	1.86	0.56
1:C:179:ALA:HB1	1:C:267:ILE:HD13	1.85	0.56
1:A:423:ARG:HG2	1:A:461:ILE:HD11	1.86	0.56
1:A:136:ILE:HD13	2:E:190:THR:HG23	1.87	0.56
1:B:344:SER:HA	7:F:600:ANP:O1G	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:253:LEU:HD23	2:F:296:ILE:HG23	1.87	0.56
3:G:37:GLU:HB3	3:G:218:LYS:HE3	1.87	0.56
1:B:148:THR:HA	1:B:182:THR:HG23	1.87	0.55
2:E:203:SER:HB2	2:E:420:VAL:HG13	1.87	0.55
1:C:183:ILE:HD11	1:C:267:ILE:HD11	1.88	0.55
2:F:244:ARG:O	2:F:248:GLY:HA2	2.06	0.55
4:H:58:LEU:HD13	4:H:77:VAL:HG11	1.87	0.55
1:B:422:VAL:O	1:B:426:GLU:HG2	2.07	0.55
1:A:169:GLY:N	1:A:175:LYS:HD3	2.22	0.55
2:F:163:THR:O	2:F:167:MET:HG2	2.06	0.54
2:F:49:VAL:HA	2:F:60:THR:HG22	1.88	0.54
3:G:20:THR:HG21	3:G:235:ALA:HB3	1.89	0.54
2:D:162:LYS:HE3	2:D:257:ASN:ND2	2.23	0.54
1:B:294:TYR:HD1	1:B:298:VAL:HG11	1.72	0.54
1:B:30:ARG:HB3	1:B:42:HIS:HB3	1.90	0.54
1:B:312:MET:HG2	1:B:319:GLY:O	2.07	0.54
1:B:241:PRO:HA	1:B:244:TYR:HB3	1.90	0.53
1:C:386:VAL:O	1:C:449:VAL:HG21	2.08	0.53
1:A:48:GLN:HB3	2:E:68:GLY:HA2	1.89	0.53
1:B:180:ILE:HG13	1:B:216:LEU:HD11	1.90	0.53
1:B:295:PRO:HG2	1:B:298:VAL:HG13	1.91	0.53
4:H:35:GLN:HG3	4:H:48:LEU:HG	1.91	0.53
1:A:148:THR:HA	1:A:182:THR:HG23	1.89	0.53
1:C:211:SER:O	1:C:215:GLN:HG2	2.08	0.53
1:A:110:ALA:HB2	1:A:234:ALA:HB2	1.90	0.52
2:D:49:VAL:HA	2:D:60:THR:HG22	1.91	0.52
1:B:47:VAL:HG12	1:B:90:ARG:HG2	1.89	0.52
2:D:89:GLU:HG3	2:D:110:THR:HA	1.90	0.52
2:D:404:VAL:O	2:D:408:ARG:HG3	2.09	0.52
1:C:183:ILE:HD12	1:C:201:CYS:SG	2.49	0.52
1:C:159:ILE:HD13	1:C:324:LEU:HD21	1.92	0.52
3:G:20:THR:HG23	3:G:232:MET:HE3	1.92	0.52
3:G:2:THR:O	3:G:6:ILE:HG12	2.09	0.51
1:B:179:ALA:HB1	1:B:267:ILE:HD13	1.92	0.51
1:B:423:ARG:HH21	1:B:458:PRO:HG3	1.74	0.51
2:E:49:VAL:HA	2:E:60:THR:HG22	1.92	0.51
2:D:97:VAL:HG21	2:D:228:ALA:HB1	1.92	0.51
2:D:196:LEU:HG	2:D:200:MET:HE2	1.92	0.51
1:C:272:SER:O	1:C:276:VAL:HG23	2.11	0.51
2:E:139:VAL:HG22	2:E:414:LEU:HB3	1.93	0.51
2:F:63:MET:HE1	2:F:231:ARG:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:164:VAL:N	7:F:600:ANP:O1A	2.43	0.51
4:H:37:ASP:HB2	4:H:64:VAL:HB	1.93	0.51
2:F:345:TYR:HA	2:F:346:PRO:C	2.35	0.50
3:G:127:THR:HB	5:I:45:VAL:HB	1.93	0.50
1:B:338:ILE:HB	1:B:339:PRO:HD3	1.92	0.50
4:H:58:LEU:HD11	4:H:92:LEU:HD11	1.93	0.50
2:F:425:THR:HG21	7:F:600:ANP:H2	1.93	0.50
1:B:357:PHE:C	1:B:359:LYS:H	2.18	0.50
2:E:27:GLU:CD	2:E:27:GLU:H	2.20	0.50
1:A:183:ILE:HD11	1:A:267:ILE:CD1	2.42	0.50
1:A:103:LEU:HD23	1:A:121:ILE:HG21	1.93	0.49
1:C:102:GLU:CD	1:C:123:SER:HA	2.38	0.49
2:E:182:VAL:HG22	2:E:216:ALA:HB3	1.94	0.49
1:A:352:LEU:HA	1:A:364:ALA:O	2.11	0.49
4:H:82:VAL:HB	4:H:92:LEU:HD13	1.94	0.49
2:E:312:VAL:HG13	2:E:322:PRO:HG3	1.94	0.49
2:F:196:LEU:HD11	2:F:200:MET:HE3	1.95	0.49
3:G:22:SER:HA	3:G:25:MET:HE3	1.93	0.49
4:H:84:VAL:HG12	4:H:90:VAL:HG22	1.95	0.49
1:C:60:LYS:O	1:C:76:PHE:HB2	2.12	0.49
1:A:353:GLU:HB2	1:A:356:LEU:HD12	1.95	0.49
1:C:457:GLU:HG2	1:C:458:PRO:HD2	1.93	0.48
1:C:66:LEU:O	2:D:15:ALA:HA	2.13	0.48
2:F:25:PHE:HB2	2:F:29:LEU:HD12	1.96	0.48
2:F:251:VAL:HB	2:F:304:ILE:HG12	1.94	0.48
2:F:293:GLN:HA	2:F:296:ILE:HD12	1.96	0.48
1:A:392:LEU:HG	1:A:396:GLN:HE21	1.77	0.48
1:B:94:ILE:O	1:B:95:VAL:C	2.56	0.48
1:B:180:ILE:O	1:B:184:ILE:HG12	2.13	0.48
2:D:370:VAL:O	2:D:374:VAL:HG23	2.13	0.48
2:E:27:GLU:CG	2:E:28:GLY:H	2.14	0.48
2:E:47:LEU:HB3	2:E:60:THR:HB	1.96	0.48
2:F:279:VAL:HG12	2:F:279:VAL:O	2.12	0.48
1:C:412:ALA:HA	1:C:415:GLN:HB2	1.95	0.48
1:A:152:ALA:HA	1:A:428:LEU:HD22	1.96	0.48
1:B:341:ASN:O	1:B:345:ILE:HG12	2.14	0.48
2:F:89:GLU:CB	2:F:110:THR:HG22	2.41	0.48
4:H:127:THR:O	4:H:131:ILE:HG12	2.14	0.48
2:D:142:LEU:HD21	2:D:374:VAL:HG21	1.95	0.47
1:B:99:VAL:CG1	1:B:100:GLY:N	2.77	0.47
2:E:32:ILE:O	2:E:33:LEU:HB2	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:VAL:HB	1:A:116:ASP:HB3	1.97	0.47
1:C:432:GLN:HE22	7:C:600:ANP:H2'	1.78	0.47
1:B:303:SER:HB2	2:F:222:MET:HB2	1.96	0.47
2:E:86:VAL:HG12	2:E:239:VAL:HA	1.96	0.47
2:F:15:ALA:HB3	2:F:22:ASP:HB2	1.96	0.47
2:D:139:VAL:HG21	2:D:346:PRO:HG2	1.95	0.47
1:B:202:ILE:HG13	1:B:230:ILE:HB	1.95	0.47
1:C:270:ASP:HB2	1:C:326:VAL:O	2.15	0.47
2:D:101:PRO:HB3	2:D:108:ILE:HD11	1.97	0.47
3:G:71:VAL:HA	3:G:108:VAL:CG1	2.44	0.47
3:G:136:PRO:O	3:G:218:LYS:NZ	2.48	0.47
3:G:130:GLU:HB3	5:I:41:THR:HG22	1.97	0.47
1:B:48:GLN:HB2	1:B:51:GLU:HB2	1.97	0.47
1:B:97:VAL:HG11	1:B:249:SER:HB3	1.97	0.47
1:B:175:LYS:HG2	1:B:352:LEU:HD12	1.97	0.47
2:D:87:GLY:HA2	2:D:242:TYR:CE1	2.50	0.47
1:C:251:CYS:O	1:C:255:GLU:HG3	2.14	0.46
4:H:23:PRO:HD3	4:H:94:ALA:O	2.15	0.46
5:I:4:TRP:H	5:I:4:TRP:CD1	2.32	0.46
1:C:162:GLY:HA3	1:C:311:LYS:HB2	1.97	0.46
6:O:39:ASN:CB	6:O:40:PRO:HA	2.45	0.46
1:B:148:THR:HG21	1:B:153:VAL:HG11	1.96	0.46
1:B:414:THR:HA	1:B:417:LEU:HD12	1.98	0.46
1:C:190:ASN:HA	1:C:198:LYS:HD3	1.97	0.46
2:F:258:ILE:HD11	2:F:292:MET:HE1	1.98	0.46
2:F:290:GLY:HA2	2:F:328:HIS:CE1	2.48	0.46
1:B:99:VAL:CG1	1:B:100:GLY:H	2.27	0.46
2:F:164:VAL:HG23	7:F:600:ANP:O1A	2.16	0.46
2:F:256:ASP:HA	2:F:257:ASN:HA	1.69	0.46
3:G:178:ILE:HD12	3:G:212:ILE:HG21	1.97	0.46
1:A:362:ARG:HH22	2:D:372:ARG:HD3	1.81	0.46
1:A:202:ILE:HG12	1:A:230:ILE:HB	1.98	0.46
1:A:457:GLU:HG2	1:A:458:PRO:HD2	1.97	0.46
2:F:61:ILE:HD11	2:F:272:LEU:HD11	1.97	0.46
3:G:164:ARG:HH21	3:G:166:ARG:HD3	1.80	0.46
2:E:259:PHE:CE1	2:E:313:PRO:HG3	2.49	0.46
1:A:98:PRO:HA	1:A:126:ARG:HG2	1.97	0.46
2:E:251:VAL:HB	2:E:304:ILE:HG12	1.98	0.46
2:E:264:ALA:O	2:E:268:VAL:HG23	2.16	0.46
3:G:113:ARG:O	3:G:117:HIS:HB2	2.16	0.46
1:A:80:LYS:HA	2:D:32:ILE:HB	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:VAL:HG12	1:A:378:ALA:HB2	1.98	0.46
1:C:289:PRO:HD2	3:G:268:GLY:HA2	1.97	0.46
1:B:225:ALA:HA	1:B:228:TYR:CE2	2.51	0.45
2:E:326:PHE:HA	2:E:329:LEU:HD12	1.97	0.45
2:E:357:ILE:HG22	2:E:362:ILE:HD13	1.97	0.45
2:F:275:ILE:HG22	2:F:276:PRO:HD2	1.98	0.45
1:A:288:PRO:HB2	2:E:270:ALA:O	2.16	0.45
1:A:362:ARG:HA	1:A:363:PRO:C	2.41	0.45
1:B:209:LYS:HD2	2:E:328:HIS:HA	1.99	0.45
2:D:89:GLU:HG3	2:D:110:THR:HG22	1.98	0.45
2:E:296:ILE:HG21	2:E:306:SER:HB2	1.98	0.45
3:G:161:ILE:HG12	3:G:175:GLU:HG2	1.96	0.45
2:F:468:ALA:O	2:F:472:LYS:HG2	2.17	0.45
2:D:345:TYR:HA	2:D:346:PRO:C	2.41	0.45
5:I:37:THR:HG22	5:I:38:SER:H	1.81	0.45
1:A:32:LEU:HG	1:A:42:HIS:HB2	1.99	0.45
2:F:258:ILE:HG21	2:F:310:ILE:HG12	1.99	0.45
2:D:256:ASP:HA	2:D:257:ASN:HA	1.70	0.45
1:B:187:LYS:HE3	1:B:224:ASP:OD2	2.16	0.45
2:F:374:VAL:HG23	2:F:445:LEU:HD11	1.99	0.45
2:E:10:THR:HG21	2:E:74:LYS:HD2	1.99	0.45
2:E:81:PRO:HB2	2:E:115:ALA:HB1	1.99	0.45
2:E:218:VAL:HG21	2:E:236:GLY:HA2	1.98	0.45
2:F:282:GLN:H	2:F:282:GLN:NE2	1.98	0.45
2:D:269:SER:O	2:D:274:ARG:HD3	2.17	0.44
2:F:29:LEU:HD12	2:F:30:PRO:HD2	1.99	0.44
1:C:137:ILE:CD1	2:D:103:ASP:HA	2.47	0.44
1:A:29:GLY:HA3	1:A:42:HIS:O	2.18	0.44
1:A:99:VAL:CG2	1:A:253:MET:HA	2.38	0.44
1:A:104:LEU:HD23	1:A:230:ILE:HG13	1.99	0.44
2:D:163:THR:O	2:D:166:ILE:HG22	2.18	0.44
2:D:239:VAL:O	2:D:243:PHE:HD2	1.99	0.44
2:E:321:ALA:HB3	2:E:322:PRO:CD	2.47	0.44
2:F:51:GLN:HG3	2:F:59:ARG:HB3	2.00	0.44
3:G:137:THR:HG22	3:G:138:PHE:N	2.33	0.44
3:G:193:TYR:HA	4:H:53:PRO:HB2	1.99	0.44
1:B:163:GLN:HG2	1:B:164:ARG:N	2.32	0.44
3:G:190:MET:HE3	3:G:190:MET:HA	1.99	0.44
1:C:336:ALA:HB3	1:C:339:PRO:HG2	1.99	0.44
2:E:36:LEU:HB2	2:E:47:LEU:HB2	2.00	0.44
2:E:460:VAL:HG21	2:E:466:ALA:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ILE:HD11	1:A:267:ILE:HD12	2.00	0.44
1:C:34:ILE:HD12	1:C:39:ALA:HB2	1.99	0.44
1:C:148:THR:HA	1:C:182:THR:HG23	2.00	0.44
2:E:345:TYR:HA	2:E:346:PRO:C	2.42	0.44
4:H:33:VAL:HG13	4:H:66:HIS:O	2.18	0.44
5:I:5:ARG:O	5:I:6:GLN:CB	2.65	0.44
1:C:338:ILE:HD13	1:C:338:ILE:HA	1.90	0.43
2:D:409:LYS:HG2	2:D:454:GLU:HA	1.99	0.43
2:E:244:ARG:HD3	2:E:304:ILE:HG13	2.00	0.43
2:E:280:GLY:HA2	3:G:262:LEU:CD2	2.48	0.43
2:E:358:MET:HE3	2:E:358:MET:HA	2.00	0.43
2:F:142:LEU:O	2:F:358:MET:HE1	2.18	0.43
1:B:185:ASN:O	1:B:188:ARG:HG2	2.18	0.43
1:C:288:PRO:HB2	2:D:270:ALA:HB1	1.99	0.43
1:C:297:ASP:HB2	1:C:300:TYR:HB3	2.00	0.43
2:D:312:VAL:HA	2:D:313:PRO:HD3	1.81	0.43
1:A:172:GLN:HA	7:A:600:ANP:N3B	2.33	0.43
1:A:383:MET:HE3	1:A:386:VAL:HG23	1.99	0.43
1:C:110:ALA:HB2	1:C:234:ALA:HB2	1.99	0.43
2:D:38:VAL:HA	2:D:75:VAL:HG22	1.99	0.43
1:A:196:LYS:H	1:A:196:LYS:HG3	1.65	0.43
1:A:496:LYS:O	1:A:499:GLU:HG2	2.18	0.43
2:E:404:VAL:O	2:E:408:ARG:HG3	2.19	0.43
2:F:231:ARG:HA	2:F:234:LEU:HD12	1.99	0.43
1:C:271:LEU:HD13	1:C:302:HIS:CD2	2.54	0.43
1:C:283:LEU:HD21	1:C:289:PRO:HB3	1.99	0.43
2:D:158:ALA:HA	7:D:600:ANP:O3G	2.18	0.43
4:H:138:ASN:HA	4:H:141:LEU:HD12	2.00	0.43
1:A:362:ARG:HH22	2:D:372:ARG:CG	2.32	0.43
2:D:89:GLU:H	2:D:89:GLU:HG2	1.46	0.43
2:E:160:VAL:HG21	2:E:335:LEU:HB3	2.00	0.43
4:H:78:SER:HB3	5:I:22:VAL:HG21	2.01	0.43
1:B:164:ARG:HG3	1:B:323:ALA:HB3	1.99	0.43
1:C:373:ARG:HG2	7:D:600:ANP:O3'	2.19	0.43
2:D:281:TYR:HB3	2:D:285:LEU:HD12	2.00	0.43
1:C:180:ILE:HD11	1:C:216:LEU:HD12	2.01	0.43
1:C:311:LYS:NZ	1:C:318:GLY:O	2.52	0.43
1:A:362:ARG:NH2	2:D:372:ARG:HD3	2.34	0.43
1:B:270:ASP:HB3	1:B:326:VAL:HB	2.00	0.43
1:B:288:PRO:HA	1:B:289:PRO:HD2	1.94	0.43
1:C:399:GLU:HG3	2:D:342:LEU:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:256:ASP:HA	2:E:257:ASN:HA	1.62	0.43
1:C:269:ASP:HA	1:C:270:ASP:HA	1.77	0.42
2:F:258:ILE:CG2	2:F:310:ILE:HG12	2.49	0.42
2:F:422:GLU:C	2:F:424:PHE:H	2.27	0.42
2:D:266:SER:HB3	2:D:282:GLN:HE22	1.84	0.42
1:A:48:GLN:HA	2:E:69:LEU:O	2.20	0.42
1:B:440:GLU:CD	1:B:473:ILE:HD11	2.43	0.42
1:B:48:GLN:HB3	2:F:68:GLY:HA2	2.01	0.42
1:B:180:ILE:HD11	1:B:216:LEU:HD21	2.01	0.42
1:B:451:GLY:C	1:B:453:LEU:H	2.27	0.42
2:E:395:GLU:HG3	3:G:36:ARG:NH2	2.33	0.42
2:F:163:THR:HA	2:F:166:ILE:HG22	2.02	0.42
1:C:286:ARG:HA	2:F:275:ILE:HD11	2.02	0.42
1:C:382:ALA:HB1	1:C:442:VAL:HG11	2.01	0.42
1:A:62:MET:O	1:A:73:VAL:HA	2.19	0.42
1:C:177:SER:OG	7:C:600:ANP:H8	2.20	0.42
1:C:418:LEU:O	1:C:422:VAL:HG23	2.19	0.42
3:G:31:TYR:CD1	3:G:226:SER:HB3	2.55	0.42
4:H:47:ILE:HG23	4:H:51:HIS:CB	2.50	0.42
4:H:47:ILE:HG12	4:H:84:VAL:HG11	2.00	0.42
2:D:36:LEU:HB2	2:D:47:LEU:HB2	2.02	0.42
2:F:410:ILE:HG13	2:F:444:ILE:HG21	2.01	0.42
4:H:65:VAL:O	4:H:72:THR:HG23	2.19	0.42
2:D:164:VAL:HG23	7:D:600:ANP:O1A	2.20	0.42
2:E:457:PHE:O	2:E:460:VAL:HG22	2.19	0.42
3:G:76:GLY:HA2	3:G:82:HIS:HE1	1.85	0.41
4:H:132:GLN:HA	4:H:135:ILE:HD12	2.01	0.41
2:F:151:LYS:HD3	2:F:328:HIS:O	2.20	0.41
3:G:42:ARG:NH2	3:G:165:PHE:O	2.53	0.41
1:A:403:PHE:CE1	3:G:19:ILE:HA	2.43	0.41
1:C:265:LEU:HD11	1:C:324:LEU:HG	2.02	0.41
2:D:66:THR:HB	2:D:69:LEU:HD12	2.02	0.41
2:D:396:LEU:HD22	2:D:400:ASP:HB3	2.03	0.41
2:F:189:ARG:O	2:F:192:GLU:HB2	2.20	0.41
1:B:176:THR:OG1	7:B:600:ANP:O2B	2.37	0.41
1:B:210:ARG:H	2:E:294:GLU:CD	2.28	0.41
2:F:153:GLY:HA3	2:F:329:LEU:HD13	2.03	0.41
1:B:68:PRO:HD3	2:F:15:ALA:HB2	2.02	0.41
1:B:397:TYR:CG	1:B:421:GLY:HA3	2.54	0.41
2:E:275:ILE:O	2:E:283:PRO:HG3	2.20	0.41
1:B:165:GLU:O	1:B:325:PRO:HD2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:187:GLY:O	2:D:222:MET:HB3	2.20	0.41
2:F:336:SER:HB3	2:F:339:ILE:HG12	2.02	0.41
2:F:440:GLY:O	2:F:444:ILE:HD12	2.21	0.41
1:B:334:VAL:HG11	1:B:351:PHE:CE2	2.56	0.41
1:C:352:LEU:HA	1:C:364:ALA:O	2.21	0.41
1:C:376:SER:HB2	1:C:384:LYS:HG3	2.01	0.41
2:D:174:ALA:O	2:D:177:HIS:HB3	2.21	0.41
2:E:168:GLU:HG2	2:E:418:PHE:CD1	2.56	0.41
1:B:26:GLU:HA	1:B:45:ARG:HB2	2.03	0.41
1:B:45:ARG:HA	1:B:45:ARG:HD3	1.88	0.41
3:G:108:VAL:HG22	3:G:131:VAL:HG21	2.03	0.41
1:C:353:GLU:HB2	1:C:356:LEU:HD12	2.01	0.41
1:C:410:LEU:HB3	1:C:414:THR:HG23	2.03	0.41
2:D:156:GLY:H	2:D:312:VAL:HG23	1.86	0.41
2:D:247:GLU:O	2:D:249:GLN:N	2.52	0.41
2:D:461:GLY:HA3	2:D:462:PRO:HD2	1.95	0.41
2:E:52:HIS:CD2	2:E:58:VAL:HG12	2.56	0.41
2:F:269:SER:HB2	2:F:282:GLN:HB3	2.01	0.41
2:F:366:GLU:O	2:F:370:VAL:HG23	2.20	0.41
2:F:408:ARG:O	2:F:412:ARG:NH1	2.54	0.41
1:B:97:VAL:HG12	1:B:111:LEU:O	2.21	0.41
1:B:159:ILE:HD13	1:B:324:LEU:HD21	2.02	0.41
1:B:210:ARG:HA	1:B:235:THR:HG21	2.02	0.41
2:D:247:GLU:C	2:D:249:GLN:H	2.29	0.41
2:E:149:GLY:HA2	2:E:304:ILE:O	2.21	0.41
2:E:275:ILE:HA	2:E:276:PRO:HD3	1.96	0.41
3:G:70:GLY:O	3:G:108:VAL:HG12	2.21	0.41
1:A:269:ASP:HA	1:A:270:ASP:HA	1.77	0.40
1:B:376:SER:C	1:B:378:ALA:H	2.28	0.40
1:C:99:VAL:HG21	1:C:127:ARG:CB	2.49	0.40
2:E:139:VAL:HG11	2:E:348:VAL:HB	2.03	0.40
2:F:143:LEU:HD22	2:F:375:GLN:HG3	2.03	0.40
3:G:86:ALA:HA	3:G:89:MET:HE3	2.03	0.40
2:D:222:MET:HA	2:D:229:ARG:HD3	2.04	0.40
2:D:235:THR:O	2:D:239:VAL:HG23	2.22	0.40
2:E:87:GLY:HA2	2:E:242:TYR:CE2	2.57	0.40
2:F:90:THR:HB	2:F:95:MET:HE1	2.02	0.40
1:B:300:TYR:HA	1:B:303:SER:OG	2.21	0.40
1:B:343:ILE:HD13	1:B:349:GLN:HG2	2.02	0.40
3:G:181:LEU:HD23	3:G:184:ILE:HB	2.03	0.40
3:G:228:ARG:O	3:G:232:MET:HG2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:237:LYS:HA	3:G:240:SER:HB2	2.03	0.40
1:B:137:ILE:N	1:B:138:PRO:CD	2.85	0.40
1:B:144:GLU:HA	1:B:145:PRO:HD3	1.89	0.40
1:C:381:ARG:H	1:C:381:ARG:HD2	1.87	0.40
2:D:63:MET:HE3	2:D:228:ALA:HA	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	490/492 (100%)	461 (94%)	25 (5%)	4 (1%)	16	49
1	B	490/492 (100%)	453 (92%)	29 (6%)	8 (2%)	7	36
1	C	490/492 (100%)	473 (96%)	16 (3%)	1 (0%)	43	74
2	D	465/467 (100%)	439 (94%)	22 (5%)	4 (1%)	14	47
2	E	464/467 (99%)	431 (93%)	28 (6%)	5 (1%)	11	43
2	F	464/467 (99%)	430 (93%)	27 (6%)	7 (2%)	8	37
3	G	257/272 (94%)	236 (92%)	20 (8%)	1 (0%)	30	62
4	H	129/131 (98%)	123 (95%)	6 (5%)	0	100	100
5	I	45/47 (96%)	40 (89%)	4 (9%)	1 (2%)	5	30
6	J	70/72 (97%)	62 (89%)	6 (9%)	2 (3%)	3	26
6	K	70/72 (97%)	64 (91%)	5 (7%)	1 (1%)	9	38
6	L	70/72 (97%)	57 (81%)	13 (19%)	0	100	100
6	M	70/72 (97%)	64 (91%)	4 (6%)	2 (3%)	3	26
6	N	70/72 (97%)	61 (87%)	9 (13%)	0	100	100
6	O	70/72 (97%)	64 (91%)	5 (7%)	1 (1%)	9	38
6	P	70/72 (97%)	64 (91%)	5 (7%)	1 (1%)	9	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	Q	70/72 (97%)	61 (87%)	7 (10%)	2 (3%)	3	26
All	All	3854/3903 (99%)	3583 (93%)	231 (6%)	40 (1%)	12	44

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	95	VAL
1	A	375	GLY
2	F	423	VAL
3	G	51	LEU
1	A	23	VAL
1	B	400	VAL
1	B	404	ALA
1	B	406	PHE
2	D	55	GLU
2	D	277	SER
2	D	474	ALA
2	E	391	LEU
2	E	472	LYS
2	F	28	GLY
2	F	44	ARG
2	F	358	MET
6	J	45	GLN
2	F	424	PHE
6	P	40	PRO
6	Q	44	GLN
1	A	194	ASP
1	B	31	VAL
1	B	373	ARG
1	B	452	TYR
5	I	6	GLN
6	O	39	ASN
2	E	223	ASN
1	A	361	ILE
2	D	248	GLY
2	E	81	PRO
2	F	279	VAL
6	M	40	PRO
1	B	375	GLY
2	E	279	VAL
1	C	95	VAL
2	F	461	GLY

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Mol	Chain	Res	Type
6	J	40	PRO
6	M	71	ILE
6	K	40	PRO
6	Q	40	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	397/397 (100%)	390 (98%)	7 (2%)	51	69
1	B	397/397 (100%)	389 (98%)	8 (2%)	48	67
1	C	397/397 (100%)	386 (97%)	11 (3%)	38	61
2	D	377/377 (100%)	373 (99%)	4 (1%)	65	74
2	E	376/377 (100%)	364 (97%)	12 (3%)	34	59
2	F	376/377 (100%)	366 (97%)	10 (3%)	39	62
3	G	225/230 (98%)	214 (95%)	11 (5%)	22	48
4	H	104/104 (100%)	102 (98%)	2 (2%)	50	68
5	I	38/38 (100%)	36 (95%)	2 (5%)	20	47
All	All	2687/2694 (100%)	2620 (98%)	67 (2%)	42	63

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	91	THR
1	A	157	VAL
1	A	196	LYS
1	A	216	LEU
1	A	244	TYR
1	A	361	ILE
1	B	20	ASP
1	B	21	THR
1	B	28	THR

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Mol	Chain	Res	Type
1	B	374	VAL
1	B	411	ASP
1	B	488	LYS
1	B	492	GLU
1	B	509	GLU
1	C	34	ILE
1	C	74	VAL
1	C	99	VAL
1	C	177	SER
1	C	202	ILE
1	C	211	SER
1	C	216	LEU
1	C	334	VAL
1	C	367	VAL
1	C	409	ASP
1	C	479	LEU
2	D	89	GLU
2	D	139	VAL
2	D	275	ILE
2	D	282	GLN
2	E	9	THR
2	E	14	VAL
2	E	27	GLU
2	E	89	GLU
2	E	175	LYS
2	E	205	VAL
2	E	215	VAL
2	E	218	VAL
2	E	257	ASN
2	E	363	VAL
2	E	390	ILE
2	E	391	LEU
2	F	23	VAL
2	F	32	ILE
2	F	43	THR
2	F	76	LEU
2	F	251	VAL
2	F	282	GLN
2	F	399	GLU
2	F	423	VAL
2	F	435	LYS
2	F	460	VAL

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Mol	Chain	Res	Type
3	G	2	THR
3	G	7	THR
3	G	69	ILE
3	G	81	ILE
3	G	88	GLN
3	G	118	ARG
3	G	126	VAL
3	G	183	THR
3	G	184	ILE
3	G	196	ILE
3	G	200	VAL
4	H	57	VAL
4	H	118	GLU
5	I	37	THR
5	I	41	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	260	ASN
1	A	379	GLN
1	A	396	GLN
1	A	416	GLN
1	A	432	GLN
1	A	503	ASN
1	B	172	GLN
1	B	186	GLN
1	B	341	ASN
1	B	379	GLN
1	B	415	GLN
1	C	42	HIS
1	C	186	GLN
1	C	215	GLN
1	C	263	HIS
1	C	302	HIS
1	C	415	GLN
1	C	432	GLN
2	D	282	GLN
2	D	328	HIS
2	D	367	HIS
2	D	379	GLN
2	D	427	HIS

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Mol	Chain	Res	Type
2	D	443	GLN
2	E	194	ASN
2	E	198	HIS
2	E	221	GLN
2	E	263	GLN
2	E	328	HIS
2	F	51	GLN
2	F	112	GLN
2	F	221	GLN
2	F	282	GLN
2	F	328	HIS
2	F	367	HIS
2	F	411	GLN
2	F	419	GLN
2	F	443	GLN
3	G	123	GLN
3	G	234	ASN
4	H	41	GLN
4	H	85	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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	$\# Z > 2$
9	GOL	B	701	-	5,5,5	0.36	0	5,5,5	0.34	0
7	ANP	A	600	8	33,33,33	2.91	11 (33%)	45,52,52	1.83	10 (22%)
7	ANP	F	600	8	33,33,33	2.84	11 (33%)	45,52,52	1.80	10 (22%)
7	ANP	B	600	8	33,33,33	2.87	11 (33%)	45,52,52	1.78	10 (22%)
7	ANP	C	600	8	33,33,33	2.91	11 (33%)	45,52,52	1.82	10 (22%)
10	SO4	E	630	-	4,4,4	0.25	0	6,6,6	0.08	0
7	ANP	D	600	8	33,33,33	2.97	12 (36%)	45,52,52	1.90	11 (24%)

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
9	GOL	B	701	-	-	0/4/4/4	-
7	ANP	A	600	8	-	2/18/38/38	0/3/3/3
7	ANP	F	600	8	-	4/18/38/38	0/3/3/3
7	ANP	B	600	8	-	2/18/38/38	0/3/3/3
7	ANP	C	600	8	-	4/18/38/38	0/3/3/3
7	ANP	D	600	8	-	5/18/38/38	0/3/3/3

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	600	ANP	PG-O1G	9.08	1.59	1.46
7	A	600	ANP	PB-O1B	8.92	1.59	1.46
7	A	600	ANP	PG-O1G	8.80	1.59	1.46
7	B	600	ANP	PG-O1G	8.70	1.59	1.46
7	C	600	ANP	PG-O1G	8.68	1.59	1.46
7	F	600	ANP	PB-O1B	8.63	1.59	1.46
7	C	600	ANP	PB-O1B	8.47	1.59	1.46
7	D	600	ANP	PB-O1B	8.41	1.58	1.46
7	F	600	ANP	PG-O1G	8.29	1.58	1.46
7	B	600	ANP	PB-O1B	8.27	1.58	1.46
7	B	600	ANP	C4-N3	5.58	1.44	1.34
7	C	600	ANP	C4-N3	5.57	1.44	1.34
7	A	600	ANP	C4-N3	5.38	1.44	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	F	600	ANP	C4-N3	5.18	1.44	1.34
7	D	600	ANP	C4-N3	5.17	1.44	1.34
7	D	600	ANP	PG-N3B	4.74	1.75	1.63
7	D	600	ANP	PB-N3B	4.63	1.75	1.63
7	B	600	ANP	PG-N3B	4.31	1.74	1.63
7	C	600	ANP	PB-N3B	4.28	1.74	1.63
7	C	600	ANP	PG-N3B	4.25	1.74	1.63
7	F	600	ANP	PG-N3B	4.23	1.74	1.63
7	B	600	ANP	PB-N3B	4.20	1.74	1.63
7	F	600	ANP	PB-N3B	4.17	1.74	1.63
7	A	600	ANP	PB-N3B	4.13	1.74	1.63
7	A	600	ANP	PG-N3B	4.12	1.74	1.63
7	F	600	ANP	C5-C4	3.95	1.46	1.39
7	C	600	ANP	C5-C4	3.92	1.46	1.39
7	D	600	ANP	C5-C4	3.76	1.45	1.39
7	B	600	ANP	C5-C4	3.60	1.45	1.39
7	A	600	ANP	C5-C4	3.38	1.45	1.39
7	D	600	ANP	C8-N7	3.28	1.38	1.31
7	C	600	ANP	C8-N7	3.25	1.37	1.31
7	B	600	ANP	C8-N7	3.24	1.37	1.31
7	A	600	ANP	C8-N7	3.21	1.37	1.31
7	F	600	ANP	C8-N7	3.18	1.37	1.31
7	C	600	ANP	C5-C6	2.89	1.49	1.41
7	B	600	ANP	C5-C6	2.86	1.49	1.41
7	F	600	ANP	C5-C6	2.86	1.49	1.41
7	D	600	ANP	C5-C6	2.83	1.48	1.41
7	A	600	ANP	C5-C6	2.67	1.48	1.41
7	A	600	ANP	C5-N7	-2.62	1.34	1.39
7	D	600	ANP	C4-N9	-2.61	1.32	1.37
7	F	600	ANP	C5-N7	-2.49	1.34	1.39
7	B	600	ANP	C4-N9	-2.48	1.32	1.37
7	F	600	ANP	C4-N9	-2.46	1.32	1.37
7	D	600	ANP	C5-N7	-2.45	1.34	1.39
7	C	600	ANP	PA-O1A	2.44	1.59	1.50
7	C	600	ANP	C5-N7	-2.44	1.34	1.39
7	B	600	ANP	PA-O1A	2.41	1.59	1.50
7	A	600	ANP	C4-N9	-2.31	1.32	1.37
7	B	600	ANP	C5-N7	-2.28	1.34	1.39
7	D	600	ANP	PA-O1A	2.23	1.58	1.50
7	C	600	ANP	C4-N9	-2.20	1.33	1.37
7	A	600	ANP	PA-O1A	2.20	1.58	1.50
7	F	600	ANP	PA-O1A	2.11	1.58	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	600	ANP	PA-O3A	2.01	1.61	1.59

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	600	ANP	C5-C4-N9	4.53	110.74	105.81
7	D	600	ANP	C4-C5-N7	-4.44	105.51	110.58
7	F	600	ANP	C5-C4-N9	4.43	110.64	105.81
7	A	600	ANP	C4-C5-N7	-4.35	105.61	110.58
7	B	600	ANP	C4-C5-N7	-4.33	105.63	110.58
7	A	600	ANP	C5-C4-N9	4.21	110.40	105.81
7	C	600	ANP	C5-C4-N9	4.21	110.40	105.81
7	F	600	ANP	C4-C5-N7	-4.11	105.88	110.58
7	C	600	ANP	C4-C5-N7	-4.10	105.89	110.58
7	B	600	ANP	C5-C4-N9	4.10	110.28	105.81
7	F	600	ANP	C5-C4-N3	-4.08	121.09	126.72
7	A	600	ANP	C5-C4-N3	-4.07	121.11	126.72
7	B	600	ANP	C6-C5-N7	4.04	139.87	132.09
7	A	600	ANP	N3-C2-N1	-4.01	122.51	128.58
7	B	600	ANP	N3-C2-N1	-3.99	122.55	128.58
7	D	600	ANP	C6-C5-N7	3.98	139.75	132.09
7	C	600	ANP	C5-C4-N3	-3.91	121.34	126.72
7	C	600	ANP	N3-C2-N1	-3.82	122.80	128.58
7	D	600	ANP	N3-C2-N1	-3.82	122.80	128.58
7	A	600	ANP	C6-C5-N7	3.81	139.43	132.09
7	B	600	ANP	C5-C4-N3	-3.71	121.61	126.72
7	C	600	ANP	C6-C5-N7	3.68	139.19	132.09
7	D	600	ANP	C5-C4-N3	-3.60	121.76	126.72
7	F	600	ANP	C6-C5-N7	3.54	138.91	132.09
7	F	600	ANP	O1G-PG-N3B	-3.47	106.65	111.77
7	F	600	ANP	N3-C2-N1	-3.29	123.60	128.58
7	D	600	ANP	C2-N1-C6	3.19	123.97	118.73
7	C	600	ANP	O1B-PB-N3B	-3.07	107.25	111.77
7	C	600	ANP	C2-N1-C6	2.99	123.64	118.73
7	D	600	ANP	O1B-PB-N3B	-2.98	107.39	111.77
7	B	600	ANP	C2-N1-C6	2.91	123.51	118.73
7	A	600	ANP	C2-N1-C6	2.83	123.38	118.73
7	D	600	ANP	C3'-C2'-C1'	2.80	106.76	101.46
7	A	600	ANP	C2-N3-C4	2.78	118.63	111.83
7	F	600	ANP	C3'-C2'-C1'	2.77	106.70	101.46
7	B	600	ANP	C2-N3-C4	2.66	118.33	111.83
7	C	600	ANP	C2-N3-C4	2.62	118.23	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	600	ANP	C2'-C1'-N9	-2.56	106.94	113.30
7	F	600	ANP	C2-N1-C6	2.53	122.89	118.73
7	C	600	ANP	C3'-C2'-C1'	2.52	106.23	101.46
7	D	600	ANP	C2-N3-C4	2.52	117.98	111.83
7	F	600	ANP	C2-N3-C4	2.51	117.96	111.83
7	F	600	ANP	C2'-C1'-N9	-2.32	107.54	113.30
7	A	600	ANP	O2B-PB-O3A	2.28	112.27	104.64
7	A	600	ANP	C3'-C2'-C1'	2.14	105.52	101.46
7	C	600	ANP	O1G-PG-N3B	-2.08	108.71	111.77
7	B	600	ANP	C3'-C2'-C1'	2.07	105.38	101.46
7	B	600	ANP	C2'-C3'-C4'	2.07	106.60	102.61
7	D	600	ANP	O2B-PB-O3A	2.05	111.50	104.64
7	B	600	ANP	O1B-PB-N3B	-2.02	108.80	111.77
7	A	600	ANP	C2'-C3'-C4'	2.01	106.49	102.61

There are no chirality outliers.

All (17) torsion outliers are listed below:

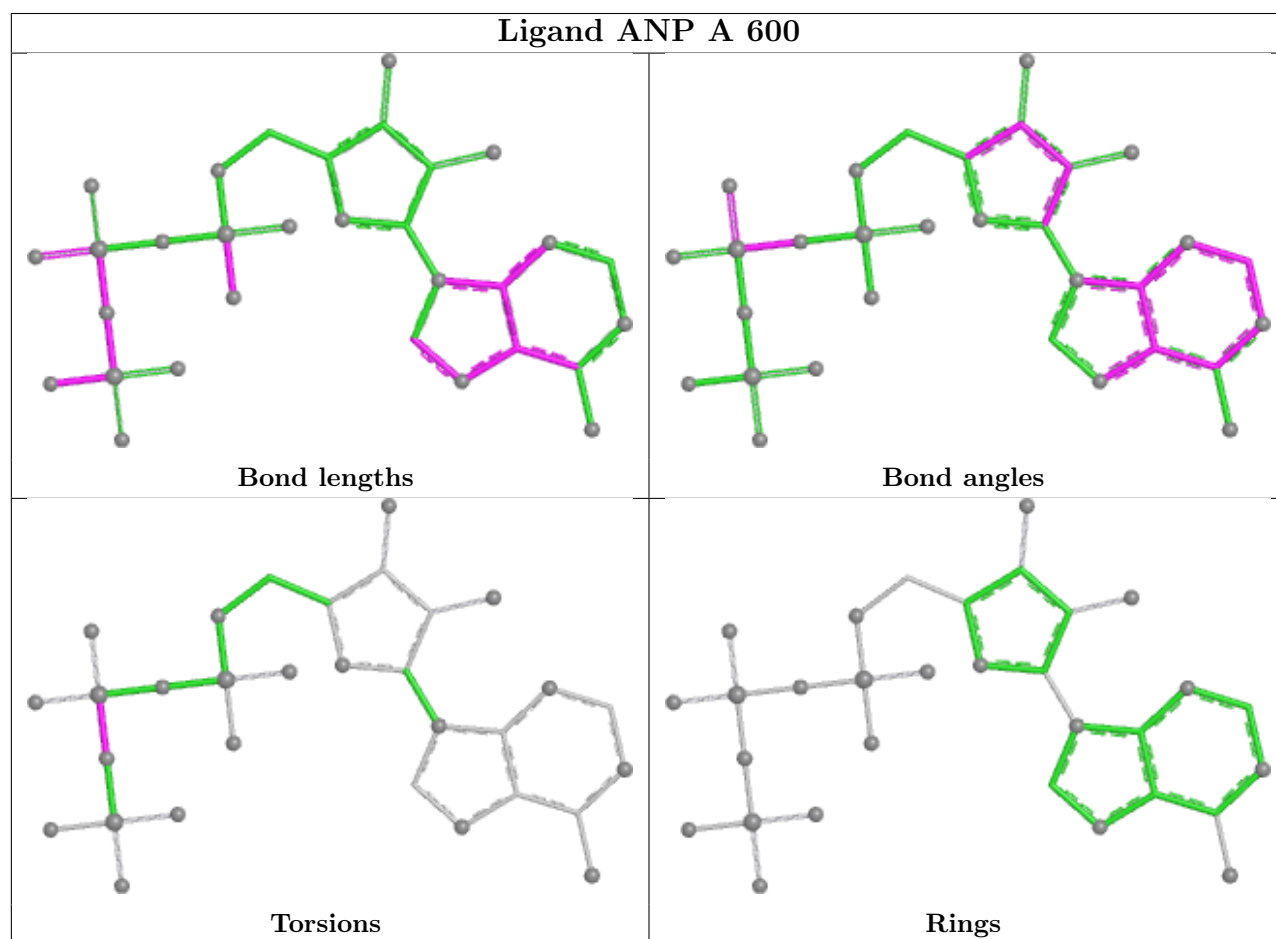
Mol	Chain	Res	Type	Atoms
7	A	600	ANP	PG-N3B-PB-O1B
7	A	600	ANP	PG-N3B-PB-O3A
7	B	600	ANP	PG-N3B-PB-O1B
7	B	600	ANP	PG-N3B-PB-O3A
7	C	600	ANP	PG-N3B-PB-O1B
7	C	600	ANP	PG-N3B-PB-O3A
7	D	600	ANP	PB-N3B-PG-O1G
7	D	600	ANP	PG-N3B-PB-O1B
7	F	600	ANP	PG-N3B-PB-O1B
7	F	600	ANP	PG-N3B-PB-O3A
7	F	600	ANP	O4'-C4'-C5'-O5'
7	D	600	ANP	O4'-C4'-C5'-O5'
7	F	600	ANP	C3'-C4'-C5'-O5'
7	C	600	ANP	O4'-C4'-C5'-O5'
7	C	600	ANP	C3'-C4'-C5'-O5'
7	D	600	ANP	C3'-C4'-C5'-O5'
7	D	600	ANP	PA-O3A-PB-O1B

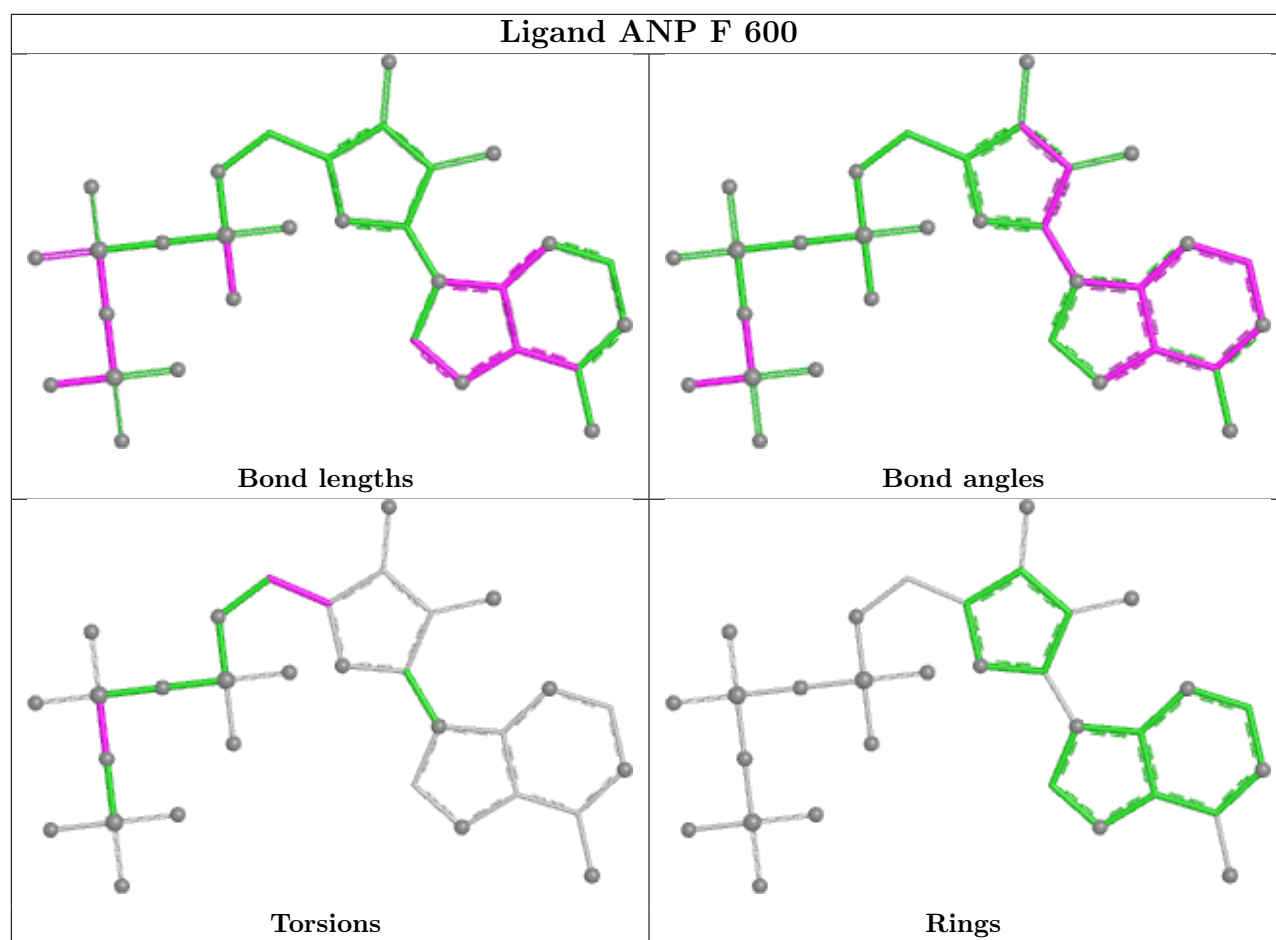
There are no ring outliers.

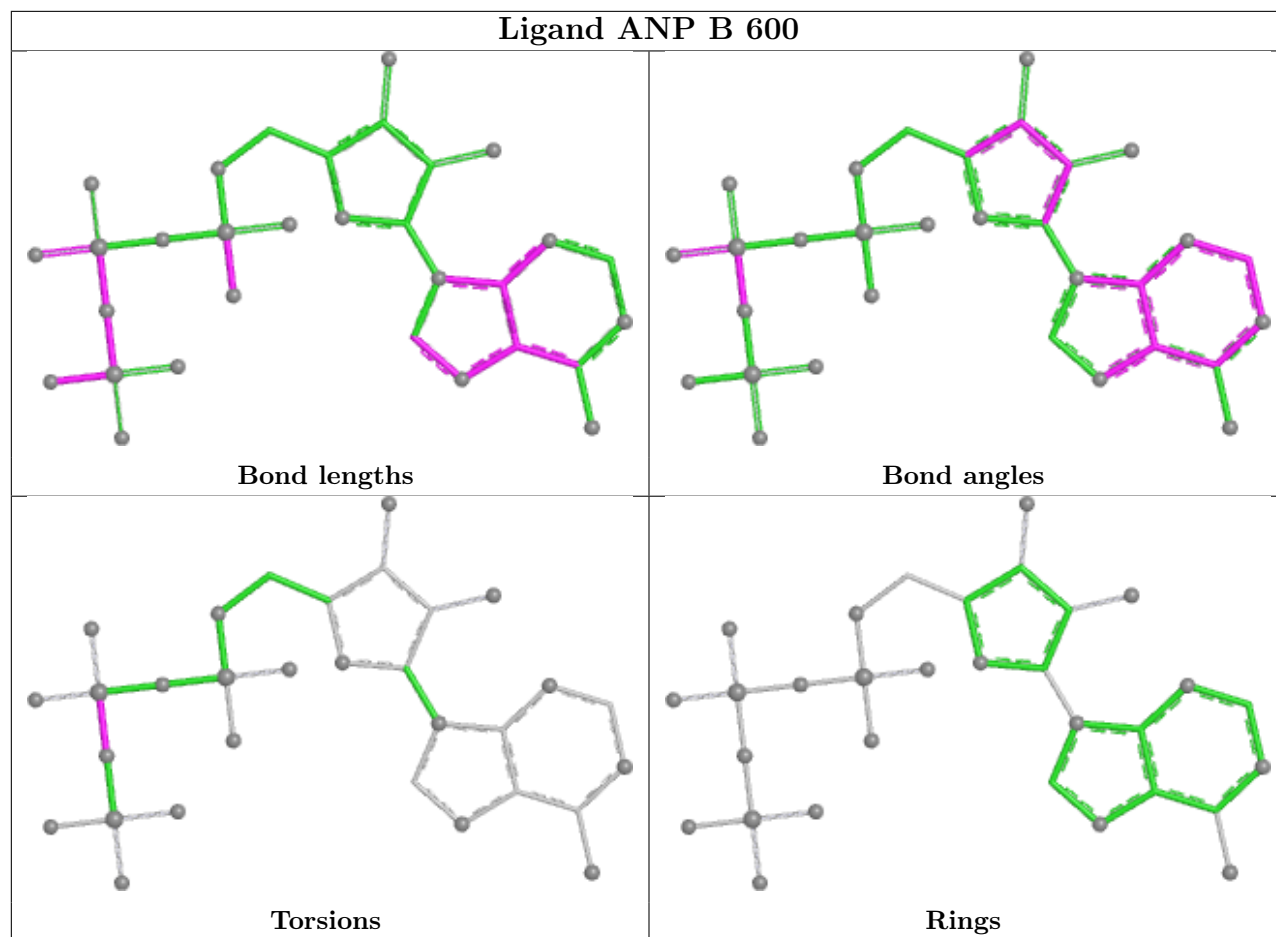
5 monomers are involved in 11 short contacts:

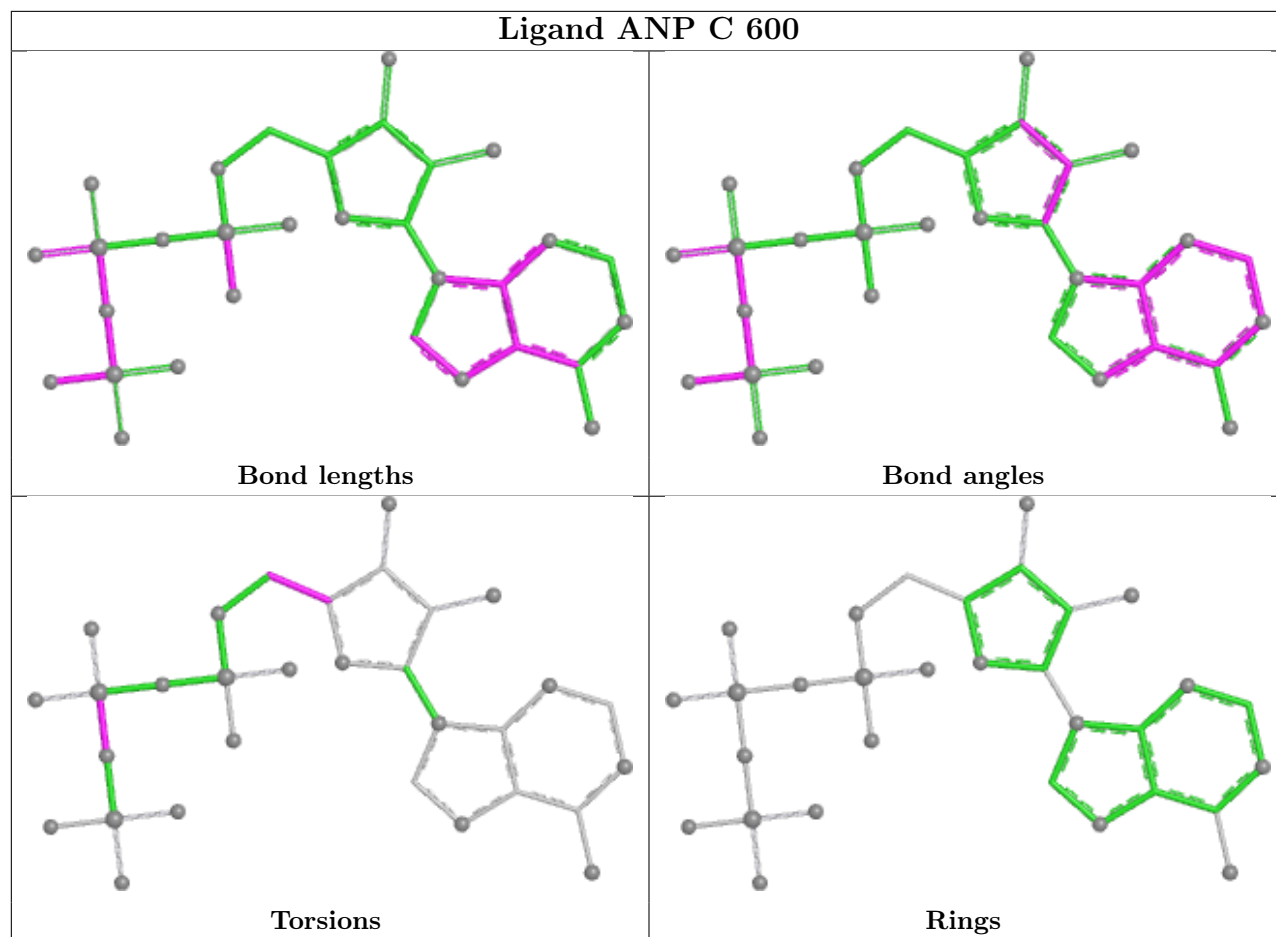
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	600	ANP	1	0
7	F	600	ANP	4	0
7	B	600	ANP	1	0
7	C	600	ANP	2	0
7	D	600	ANP	3	0

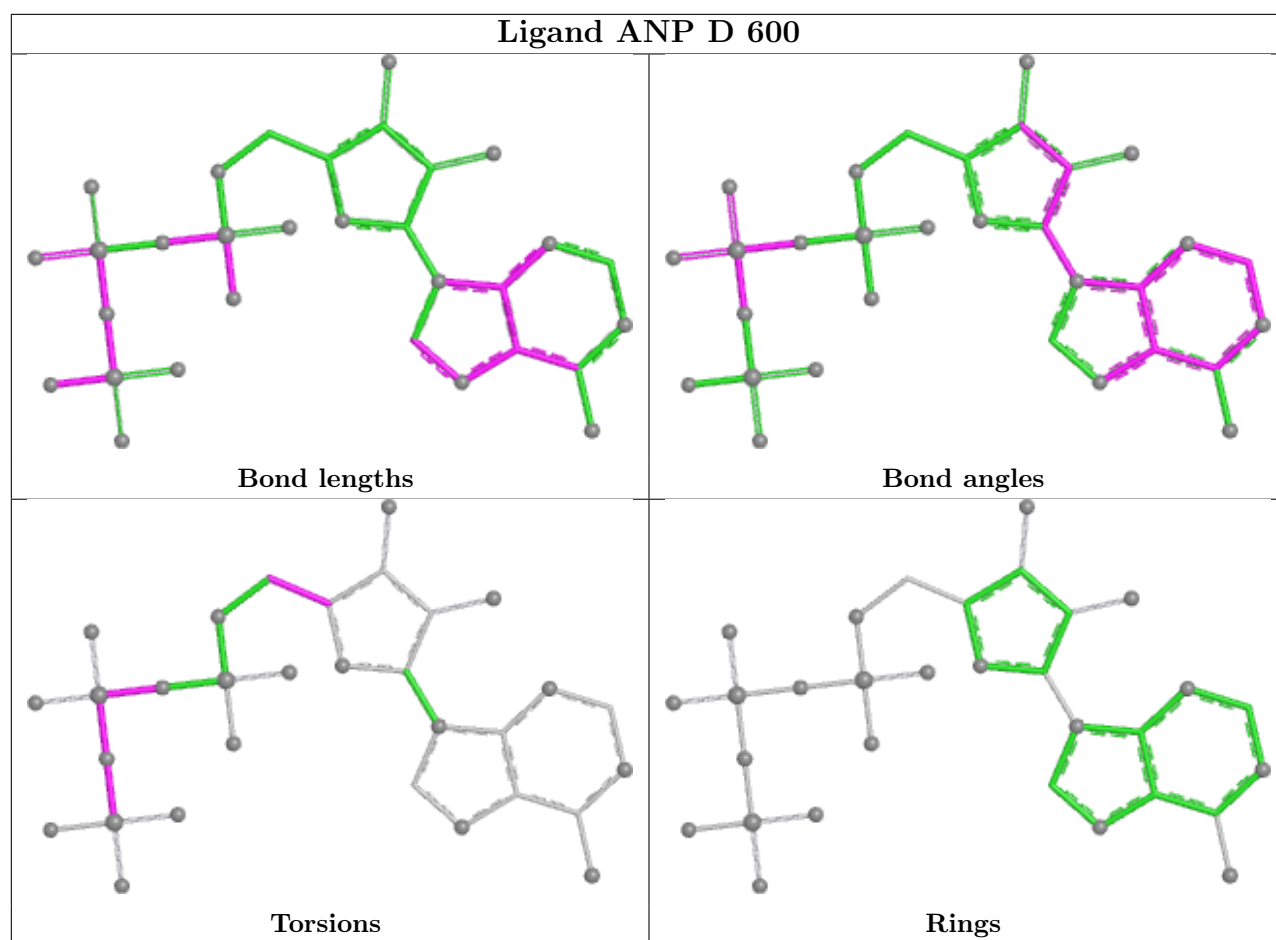
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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, 95th 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 ²)	Q<0.9
1	A	492/492 (100%)	0.10	10 (2%) 65 39	49, 67, 100, 111	0
1	B	492/492 (100%)	0.24	21 (4%) 40 21	51, 72, 110, 128	0
1	C	492/492 (100%)	0.17	12 (2%) 59 35	50, 72, 105, 107	0
2	D	467/467 (100%)	0.07	4 (0%) 81 58	49, 69, 86, 104	0
2	E	466/467 (99%)	0.18	9 (1%) 66 41	51, 74, 109, 125	0
2	F	466/467 (99%)	0.13	9 (1%) 66 41	53, 70, 87, 92	0
3	G	263/272 (96%)	0.63	17 (6%) 25 15	53, 115, 136, 141	0
4	H	131/131 (100%)	1.53	30 (22%) 2 2	138, 143, 150, 151	0
5	I	47/47 (100%)	0.97	4 (8%) 16 11	119, 125, 140, 140	0
6	J	72/72 (100%)	1.22	12 (16%) 4 4	127, 130, 135, 136	0
6	K	72/72 (100%)	1.24	14 (19%) 3 3	129, 132, 134, 134	0
6	L	72/72 (100%)	1.15	10 (13%) 6 5	129, 132, 137, 137	0
6	M	72/72 (100%)	1.08	8 (11%) 10 7	118, 127, 135, 136	0
6	N	72/72 (100%)	1.17	14 (19%) 3 3	123, 126, 136, 137	0
6	O	72/72 (100%)	1.25	15 (20%) 2 3	123, 127, 137, 137	0
6	P	72/72 (100%)	1.22	11 (15%) 5 5	128, 130, 132, 132	0
6	Q	72/72 (100%)	1.23	11 (15%) 5 5	125, 126, 132, 133	0
All	All	3892/3903 (99%)	0.39	211 (5%) 31 18	49, 76, 136, 151	0

All (211) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	86	ALA	9.1
6	P	25	ILE	5.8
4	H	55	LEU	5.6
3	G	191	SER	5.4

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Mol	Chain	Res	Type	RSRZ
6	M	19	ALA	5.4
1	B	400	VAL	5.3
3	G	180	SER	5.2
4	H	38	VAL	5.2
4	H	100	LEU	5.2
3	G	192	ILE	4.7
4	H	54	THR	4.7
6	L	3	ASP	4.5
6	L	36	TYR	4.3
3	G	196	ILE	4.3
2	E	387	ILE	4.2
4	H	81	SER	4.1
1	B	125	ALA	4.1
1	B	452	TYR	4.0
3	G	53	GLU	4.0
6	Q	16	VAL	3.9
6	J	36	TYR	3.9
6	N	11	ALA	3.8
6	Q	13	ALA	3.8
6	J	16	VAL	3.8
4	H	85	ASN	3.8
6	J	60	MET	3.8
6	K	17	GLY	3.7
4	H	45	PHE	3.6
1	A	446	TYR	3.6
6	L	19	ALA	3.6
1	B	413	ALA	3.6
6	O	34	ILE	3.5
2	E	391	LEU	3.5
1	B	447	ALA	3.5
1	B	510	ALA	3.5
6	P	17	GLY	3.5
2	E	198	HIS	3.5
6	J	37	ALA	3.4
6	O	16	VAL	3.4
4	H	75	TYR	3.4
6	N	67	VAL	3.3
6	O	10	GLY	3.3
5	I	47	VAL	3.2
6	J	15	THR	3.2
5	I	30	PHE	3.2
3	G	194	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
3	G	178	ILE	3.2
2	F	42	GLU	3.2
6	O	13	ALA	3.2
6	P	41	SER	3.2
6	P	43	LYS	3.2
6	P	16	VAL	3.2
4	H	97	ALA	3.1
6	N	13	ALA	3.1
6	K	18	VAL	3.1
6	P	29	PHE	3.1
6	Q	19	ALA	3.1
6	N	12	GLY	3.1
6	M	23	ALA	3.1
6	Q	27	THR	3.1
1	A	93	ALA	3.0
6	J	64	CYS	3.0
6	K	9	ILE	3.0
6	Q	2	ILE	3.0
6	N	27	THR	2.9
6	J	19	ALA	2.9
1	B	448	GLY	2.9
2	F	423	VAL	2.9
3	G	5	ASP	2.8
4	H	41	GLN	2.8
6	N	17	GLY	2.8
5	I	8	GLY	2.8
1	B	453	LEU	2.8
1	C	293	ALA	2.8
2	F	474	ALA	2.8
1	B	389	THR	2.8
6	O	15	THR	2.8
3	G	51	LEU	2.8
6	O	49	TYR	2.8
1	B	355	GLU	2.8
4	H	42	THR	2.8
4	H	44	ALA	2.8
6	K	3	ASP	2.7
1	C	207	GLY	2.7
1	C	427	LEU	2.7
6	N	9	ILE	2.7
3	G	187	ALA	2.7
6	Q	45	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
6	O	11	ALA	2.7
6	Q	41	SER	2.7
4	H	56	GLN	2.7
6	J	9	ILE	2.7
4	H	120	LEU	2.7
6	N	18	VAL	2.6
6	K	4	THR	2.6
4	H	52	VAL	2.6
2	F	425	THR	2.6
2	E	390	ILE	2.6
4	H	142	VAL	2.6
1	C	194	ASP	2.6
6	P	2	ILE	2.6
2	D	117	HIS	2.6
2	E	152	ILE	2.6
1	A	23	VAL	2.6
6	K	64	CYS	2.5
4	H	57	VAL	2.5
4	H	77	VAL	2.5
2	D	386	ASP	2.5
6	M	60	MET	2.5
6	N	41	SER	2.5
2	D	395	GLU	2.5
1	C	417	LEU	2.5
4	H	93	LEU	2.5
3	G	52	TYR	2.5
1	A	491	GLU	2.5
1	B	114	ALA	2.5
2	F	175	LYS	2.5
2	E	197	TYR	2.5
4	H	105	LEU	2.5
1	C	152	ALA	2.5
6	K	68	ALA	2.5
6	O	9	ILE	2.5
6	Q	34	ILE	2.5
6	L	4	THR	2.5
6	P	27	THR	2.5
6	N	16	VAL	2.5
6	K	19	ALA	2.4
1	B	113	ASN	2.4
6	J	3	ASP	2.4
1	C	23	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
6	K	43	LYS	2.4
6	N	40	PRO	2.4
3	G	197	ASP	2.4
4	H	26	VAL	2.4
6	Q	73	PHE	2.4
4	H	51	HIS	2.4
2	E	386	ASP	2.4
3	G	195	ASP	2.4
6	O	6	ALA	2.4
1	C	459	SER	2.3
6	Q	58	GLU	2.3
3	G	112	ILE	2.3
4	H	43	GLY	2.3
1	B	196	LYS	2.3
6	J	43	LYS	2.3
3	G	1	ALA	2.3
1	A	137	ILE	2.3
6	K	62	LEU	2.3
1	A	91	THR	2.3
6	N	19	ALA	2.3
6	O	17	GLY	2.3
6	J	49	TYR	2.3
1	B	371	VAL	2.3
1	A	21	THR	2.3
6	J	21	SER	2.3
6	L	16	VAL	2.3
2	E	102	ILE	2.2
6	K	55	ALA	2.2
6	N	37	ALA	2.2
6	Q	55	ALA	2.2
6	M	44	GLN	2.2
6	M	38	ARG	2.2
2	F	428	LEU	2.2
1	A	495	ALA	2.2
3	G	193	TYR	2.2
4	H	103	LEU	2.2
6	O	18	VAL	2.2
6	K	2	ILE	2.2
4	H	122	ALA	2.2
1	B	23	VAL	2.2
6	P	40	PRO	2.2
2	D	184	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	25	LEU	2.2
6	O	2	ILE	2.2
6	P	18	VAL	2.1
6	M	15	THR	2.1
6	K	38	ARG	2.1
1	C	316	PHE	2.1
1	B	390	MET	2.1
6	P	60	MET	2.1
6	L	37	ALA	2.1
6	L	29	PHE	2.1
6	L	35	GLY	2.1
2	F	318	THR	2.1
6	M	37	ALA	2.1
6	O	43	LYS	2.1
6	L	18	VAL	2.1
6	M	42	LEU	2.1
2	F	43	THR	2.1
1	B	386	VAL	2.1
1	B	449	VAL	2.1
3	G	103	VAL	2.1
4	H	145	LEU	2.1
5	I	1	VAL	2.1
1	A	37	GLY	2.0
6	O	36	TYR	2.0
4	H	53	PRO	2.0
1	A	181	ASP	2.0
1	B	410	LEU	2.0
2	F	361	ASN	2.0
6	O	24	GLY	2.0
1	C	21	THR	2.0
2	E	344	ILE	2.0
6	K	27	THR	2.0
1	C	314	ASP	2.0
1	B	392	LEU	2.0
1	C	240	ALA	2.0
6	L	21	SER	2.0
4	H	84	VAL	2.0
6	N	15	THR	2.0
4	H	94	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

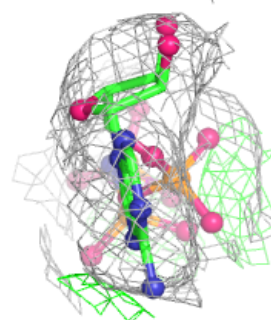
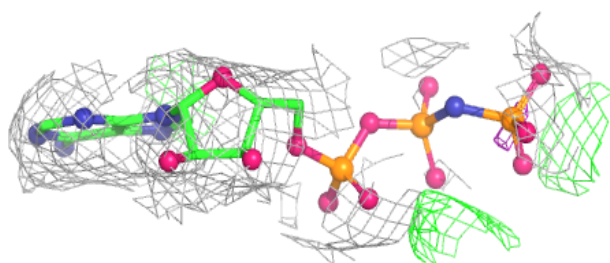
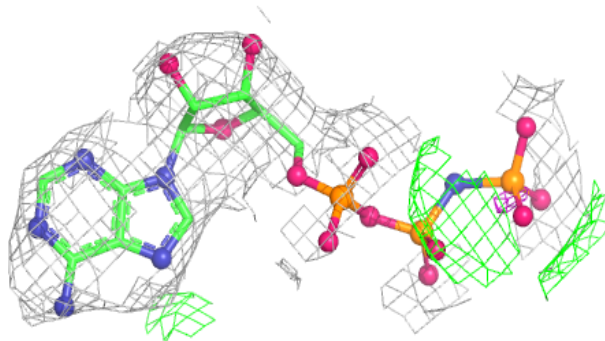
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
8	MG	D	601	1/1	0.81	0.23	62,62,62,62	0
8	MG	C	601	1/1	0.86	0.24	60,60,60,60	0
8	MG	A	601	1/1	0.87	0.25	43,43,43,43	0
8	MG	F	601	1/1	0.91	0.17	56,56,56,56	0
8	MG	B	601	1/1	0.92	0.15	55,55,55,55	0
7	ANP	D	600	31/31	0.94	0.09	62,63,67,68	0
10	SO4	E	630	5/5	0.94	0.10	71,71,71,71	0
7	ANP	F	600	31/31	0.95	0.10	55,60,61,62	0
7	ANP	C	600	31/31	0.96	0.07	59,62,64,64	0
7	ANP	A	600	31/31	0.96	0.07	41,42,43,43	0
9	GOL	B	701	6/6	0.96	0.07	42,43,43,43	0
7	ANP	B	600	31/31	0.96	0.09	54,59,61,61	0

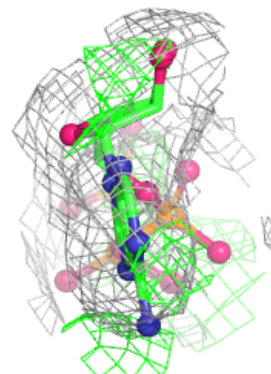
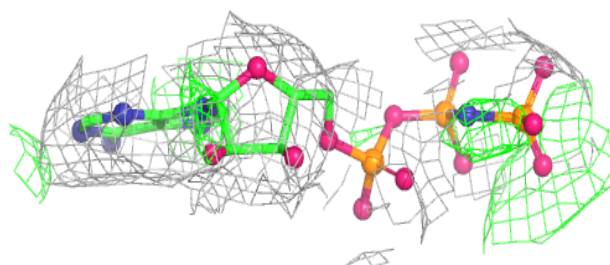
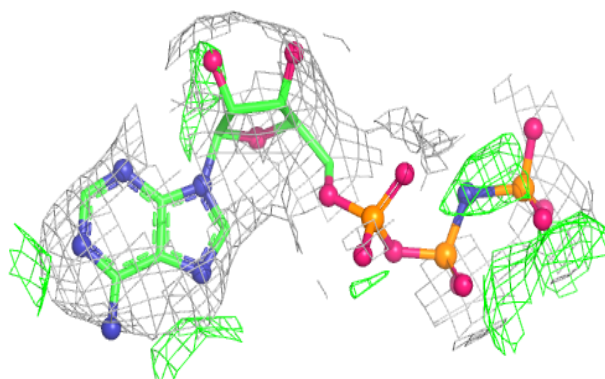
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ANP D 600:

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

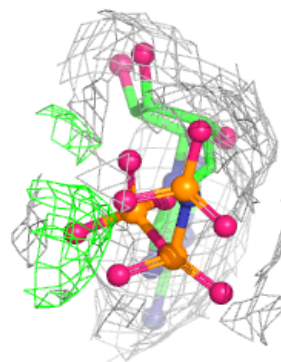
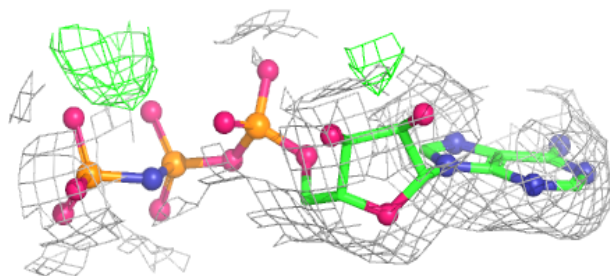
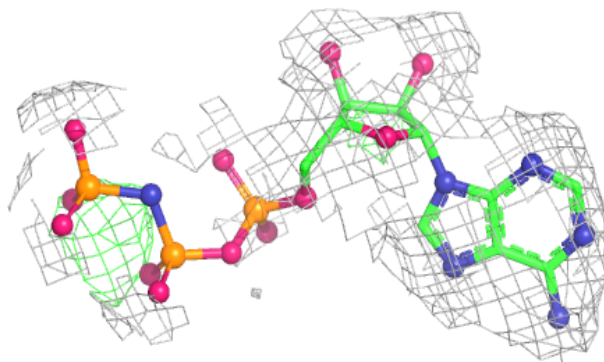
**Electron density around ANP F 600:**

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

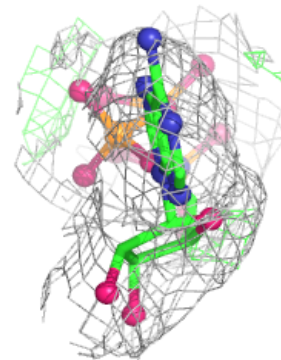
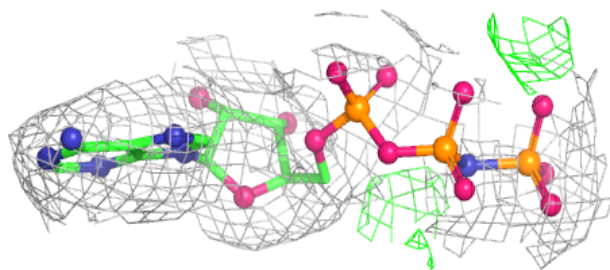
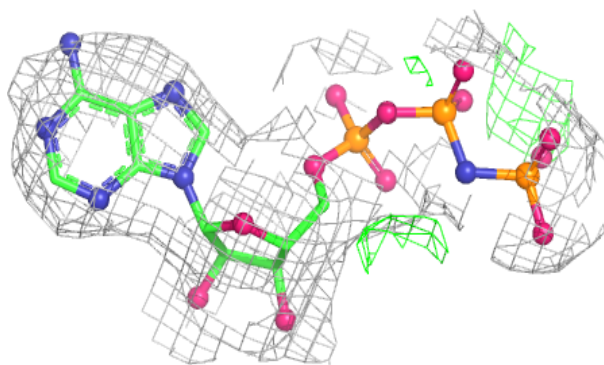


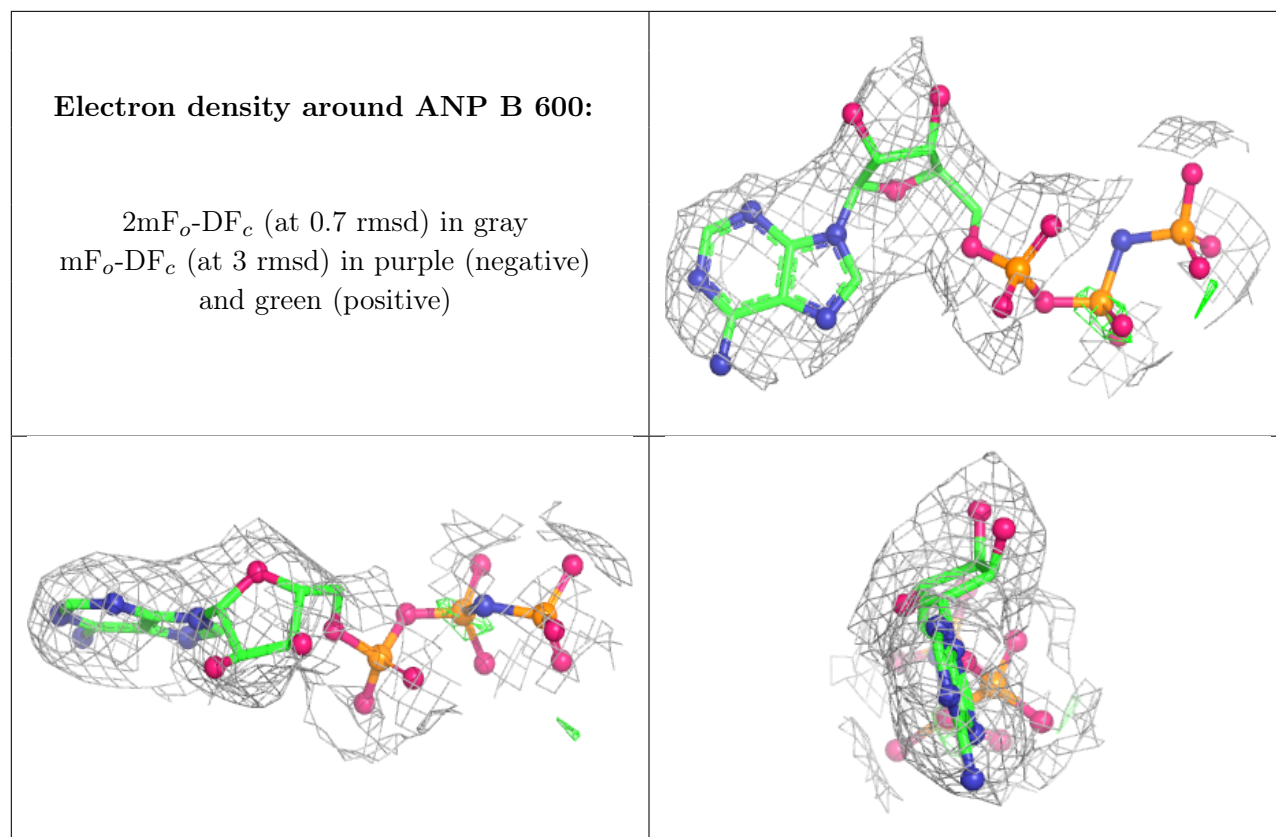
Electron density around ANP C 600:

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

**Electron density around ANP A 600:**

$2mF_o-DF_c$ (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.