



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:54 AM UTC

PDB ID : 4FI3 / pdb_00004fi3
Title : Structure of vitamin B12 transporter BtuCD-F in a nucleotide-bound state
Authors : Korkhov, V.M.; Mireku, S.A.; Locher, K.P.
Deposited on : 2012-06-07
Resolution : 3.47 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

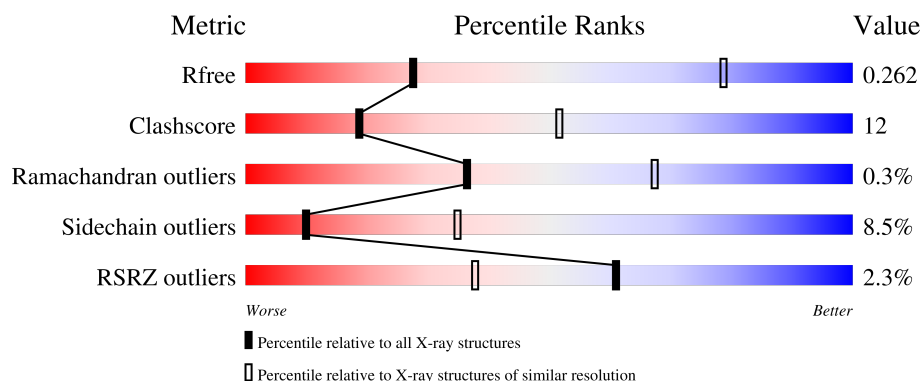
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

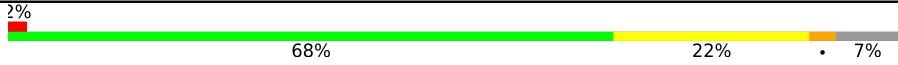
The reported resolution of this entry is 3.47 Å.

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	1070 (3.50-3.42)
Clashscore	190562	1128 (3.50-3.42)
Ramachandran outliers	187476	1101 (3.50-3.42)
Sidechain outliers	187428	1102 (3.50-3.42)
RSRZ outliers	180081	1070 (3.50-3.42)

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	349	
1	B	349	
2	C	249	
2	D	249	
3	F	255	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10635 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 Vitamin B12 import system permease protein BtuC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2441	1611	418	400	12			
1	B	324	Total	C	N	O	S	0	0	0
			2441	1611	418	400	12			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	expression tag	UNP P06609
A	-21	GLY	-	expression tag	UNP P06609
A	-20	HIS	-	expression tag	UNP P06609
A	-19	HIS	-	expression tag	UNP P06609
A	-18	HIS	-	expression tag	UNP P06609
A	-17	HIS	-	expression tag	UNP P06609
A	-16	HIS	-	expression tag	UNP P06609
A	-15	HIS	-	expression tag	UNP P06609
A	-14	HIS	-	expression tag	UNP P06609
A	-13	HIS	-	expression tag	UNP P06609
A	-12	HIS	-	expression tag	UNP P06609
A	-11	HIS	-	expression tag	UNP P06609
A	-10	SER	-	expression tag	UNP P06609
A	-9	SER	-	expression tag	UNP P06609
A	-8	GLY	-	expression tag	UNP P06609
A	-7	GLU	-	expression tag	UNP P06609
A	-6	ASN	-	expression tag	UNP P06609
A	-5	LEU	-	expression tag	UNP P06609
A	-4	TYR	-	expression tag	UNP P06609
A	-3	PHE	-	expression tag	UNP P06609
A	-2	GLN	-	expression tag	UNP P06609
A	-1	GLY	-	expression tag	UNP P06609
A	0	HIS	-	expression tag	UNP P06609
A	18	SER	CYS	engineered mutation	UNP P06609
A	32	SER	CYS	engineered mutation	UNP P06609

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Chain	Residue	Modelled	Actual	Comment	Reference
A	120	SER	CYS	engineered mutation	UNP P06609
A	156	SER	CYS	engineered mutation	UNP P06609
A	205	SER	CYS	engineered mutation	UNP P06609
A	206	SER	CYS	engineered mutation	UNP P06609
A	267	SER	CYS	engineered mutation	UNP P06609
B	-22	MET	-	expression tag	UNP P06609
B	-21	GLY	-	expression tag	UNP P06609
B	-20	HIS	-	expression tag	UNP P06609
B	-19	HIS	-	expression tag	UNP P06609
B	-18	HIS	-	expression tag	UNP P06609
B	-17	HIS	-	expression tag	UNP P06609
B	-16	HIS	-	expression tag	UNP P06609
B	-15	HIS	-	expression tag	UNP P06609
B	-14	HIS	-	expression tag	UNP P06609
B	-13	HIS	-	expression tag	UNP P06609
B	-12	HIS	-	expression tag	UNP P06609
B	-11	HIS	-	expression tag	UNP P06609
B	-10	SER	-	expression tag	UNP P06609
B	-9	SER	-	expression tag	UNP P06609
B	-8	GLY	-	expression tag	UNP P06609
B	-7	GLU	-	expression tag	UNP P06609
B	-6	ASN	-	expression tag	UNP P06609
B	-5	LEU	-	expression tag	UNP P06609
B	-4	TYR	-	expression tag	UNP P06609
B	-3	PHE	-	expression tag	UNP P06609
B	-2	GLN	-	expression tag	UNP P06609
B	-1	GLY	-	expression tag	UNP P06609
B	0	HIS	-	expression tag	UNP P06609
B	18	SER	CYS	engineered mutation	UNP P06609
B	32	SER	CYS	engineered mutation	UNP P06609
B	120	SER	CYS	engineered mutation	UNP P06609
B	156	SER	CYS	engineered mutation	UNP P06609
B	205	SER	CYS	engineered mutation	UNP P06609
B	206	SER	CYS	engineered mutation	UNP P06609
B	267	SER	CYS	engineered mutation	UNP P06609

- Molecule 2 is a protein called Vitamin B12 import ATP-binding protein BtuD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	248	Total	C	N	O	S	0	0	0
			1891	1183	351	348	9			
2	D	248	Total	C	N	O	S	0	0	0
			1891	1183	351	348	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	159	GLN	GLU	engineered mutation	UNP P06611
C	162	CYS	ASN	engineered mutation	UNP P06611
C	180	SER	CYS	engineered mutation	UNP P06611
D	159	GLN	GLU	engineered mutation	UNP P06611
D	162	CYS	ASN	engineered mutation	UNP P06611
D	180	SER	CYS	engineered mutation	UNP P06611

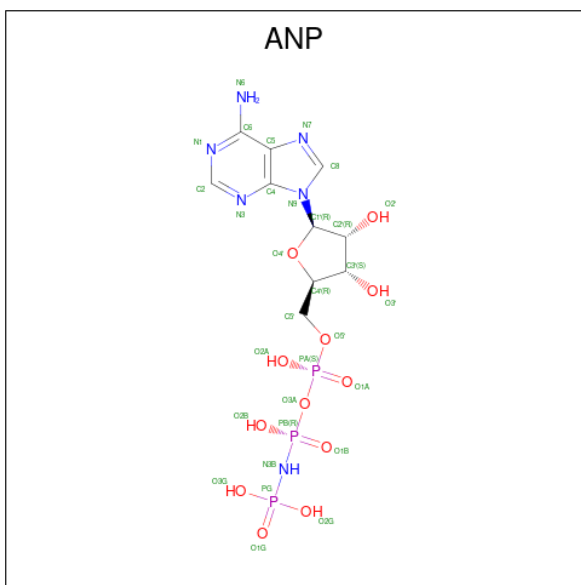
- Molecule 3 is a protein called Vitamin B12-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	245	Total	C	N	O	S	0	0	0
			1907	1216	332	355	4			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	21	MET	-	expression tag	UNP P37028
F	267	SER	-	expression tag	UNP P37028
F	268	GLY	-	expression tag	UNP P37028
F	269	SER	-	expression tag	UNP P37028
F	270	HIS	-	expression tag	UNP P37028
F	271	HIS	-	expression tag	UNP P37028
F	272	HIS	-	expression tag	UNP P37028
F	273	HIS	-	expression tag	UNP P37028
F	274	HIS	-	expression tag	UNP P37028
F	275	HIS	-	expression tag	UNP P37028

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

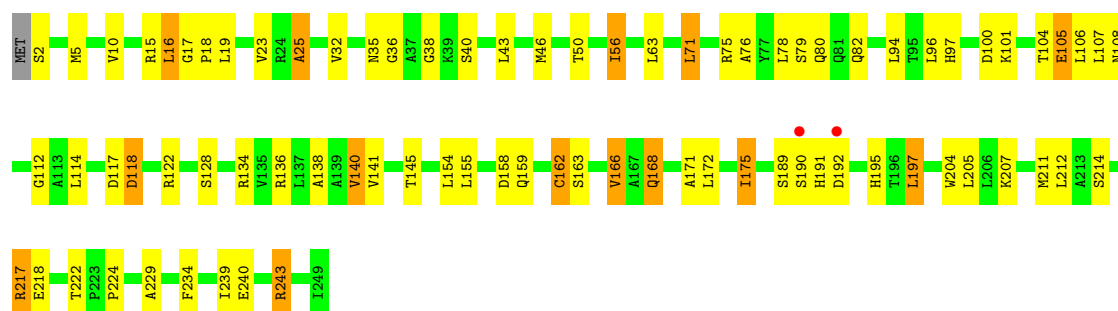
- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total Mg 1 1	0	0
5	D	1	Total Mg 1 1	0	0

i

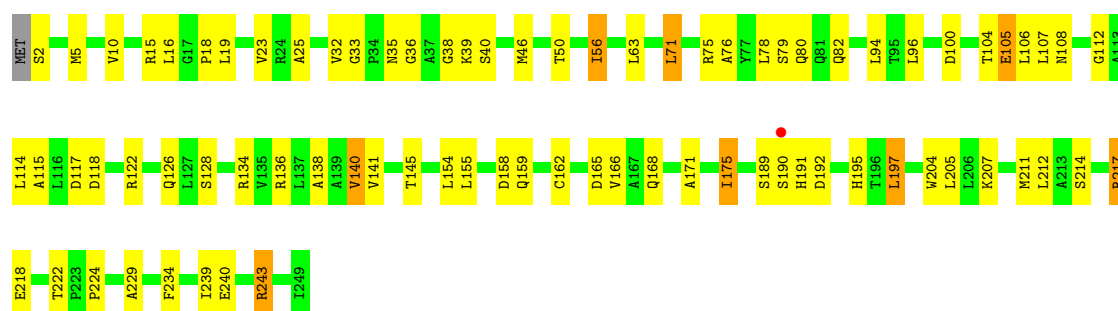
- Molecule 1: Vitamin B12 import system permease protein BtuC





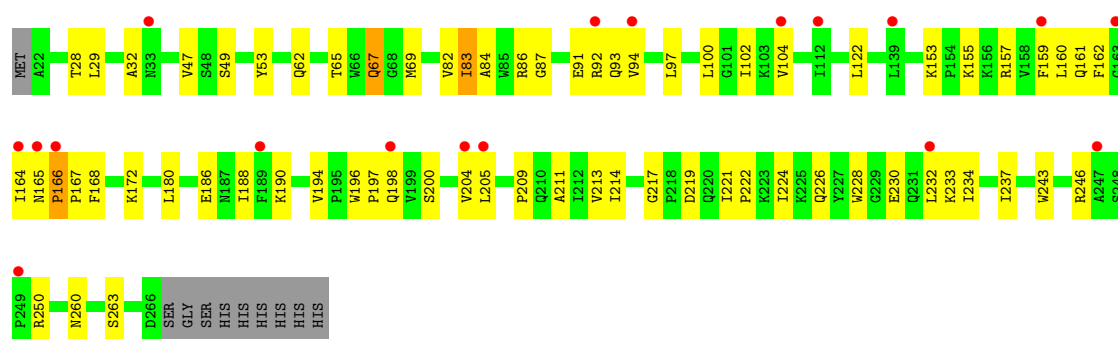
- Molecule 2: Vitamin B12 import ATP-binding protein BtuD

Chain D: 67% 29%



- Molecule 3: Vitamin B12-binding protein

Chain F: 7% 69% 26%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	127.73Å 211.99Å 179.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.47 30.00 – 3.47	Depositor EDS
% Data completeness (in resolution range)	96.3 (30.00-3.47) 96.2 (30.00-3.47)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 3.47Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.229 , 0.274 0.220 , 0.262	Depositor DCC
R_{free} test set	2007 reflections (6.54%)	wwPDB-VP
Wilson B-factor (Å ²)	118.0	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 109.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.002 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.020 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10635	wwPDB-VP
Average B, all atoms (Å ²)	169.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: ANP, MG

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	0.45	0/2495	0.84	4/3408 (0.1%)
1	B	0.43	0/2495	0.85	6/3408 (0.2%)
2	C	0.51	0/1925	0.87	3/2608 (0.1%)
2	D	0.47	0/1925	0.88	2/2608 (0.1%)
3	F	0.37	0/1951	0.78	4/2661 (0.2%)
All	All	0.45	0/10791	0.84	19/14693 (0.1%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	145	THR	CA-C-N	6.42	126.11	119.56
2	D	145	THR	C-N-CA	6.42	126.11	119.56
1	A	87	GLU	CA-C-N	5.32	125.38	119.32
1	A	87	GLU	C-N-CA	5.32	125.38	119.32
2	C	145	THR	CA-C-N	5.31	124.97	119.56
2	C	145	THR	C-N-CA	5.31	124.97	119.56
1	B	83	ASN	CA-C-N	5.21	124.87	119.56
1	B	83	ASN	C-N-CA	5.21	124.87	119.56
3	F	237	ILE	N-CA-C	5.15	112.65	107.55
1	B	87	GLU	CA-C-N	5.12	125.16	119.32
1	B	87	GLU	C-N-CA	5.12	125.16	119.32
3	F	153	LYS	CA-C-N	5.11	125.11	119.90
3	F	153	LYS	C-N-CA	5.11	125.11	119.90
2	C	25	ALA	N-CA-C	5.07	116.50	110.97
1	A	112	LEU	CA-C-N	-5.03	114.05	120.23
1	A	112	LEU	C-N-CA	-5.03	114.05	120.23
3	F	211	ALA	N-CA-C	5.01	116.46	108.79
1	B	112	LEU	CA-C-N	-5.00	114.08	120.23
1	B	112	LEU	C-N-CA	-5.00	114.08	120.23

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	2441	0	2606	65	0
1	B	2441	0	2606	63	0
2	C	1891	0	1925	53	0
2	D	1891	0	1925	47	0
3	F	1907	0	1924	53	0
4	C	31	0	13	6	0
4	D	31	0	13	5	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
All	All	10635	0	11012	257	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 12.

All (257) 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)
3:F:168:PHE:HB2	3:F:198:GLN:HE22	1.42	0.84
1:B:111:GLN:HB3	1:B:113:PRO:HD2	1.59	0.84
1:A:167:SER:HB3	1:A:171:ASP:HB2	1.61	0.82
2:C:5:MET:HB3	2:C:23:VAL:HB	1.59	0.82
1:A:111:GLN:HB3	1:A:113:PRO:HD2	1.60	0.82
1:A:98:GLY:HA3	1:A:160:MET:HE1	1.63	0.81
2:D:5:MET:HB3	2:D:23:VAL:HB	1.62	0.80
2:D:218:GLU:OE2	2:D:243:ARG:NH1	2.15	0.80
2:C:218:GLU:OE2	2:C:243:ARG:NH1	2.16	0.78
3:F:214:ILE:HD13	3:F:224:ILE:HG13	1.66	0.77
1:B:98:GLY:HA3	1:B:160:MET:HE1	1.67	0.77
1:B:261:PRO:HB2	1:B:265:ARG:HH21	1.49	0.77
1:A:261:PRO:HB2	1:A:265:ARG:HH21	1.50	0.76
1:B:129:THR:HG21	1:B:240:GLY:HA3	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:HG21	1:A:240:GLY:HA3	1.70	0.73
2:C:205:LEU:HD21	2:C:229:ALA:HB2	1.71	0.72
2:D:205:LEU:HD21	2:D:229:ALA:HB2	1.73	0.71
1:A:298:LEU:HD11	1:A:303:LEU:HB2	1.73	0.70
2:C:134:ARG:HD3	2:C:168:GLN:HG3	1.73	0.70
2:D:56:ILE:HD11	2:D:63:LEU:HD22	1.74	0.69
1:B:170:VAL:HA	3:F:87:GLY:HA2	1.74	0.69
1:B:298:LEU:HD11	1:B:303:LEU:HB2	1.73	0.69
2:C:56:ILE:HD11	2:C:63:LEU:HD22	1.76	0.68
1:B:65:LEU:HB3	1:B:197:LEU:HD23	1.78	0.66
2:C:168:GLN:OE1	2:C:168:GLN:HA	1.94	0.66
2:D:40:SER:OG	2:D:158:ASP:OD2	2.13	0.66
1:B:54:GLN:HA	1:B:188:TRP:HE1	1.62	0.65
1:A:54:GLN:HA	1:A:188:TRP:HE1	1.62	0.65
1:A:177:TYR:CD2	3:F:198:GLN:HG2	2.31	0.65
1:A:144:ARG:HH12	1:B:324:ALA:HB3	1.61	0.64
2:D:207:LYS:HD2	2:D:229:ALA:HA	1.78	0.64
1:A:65:LEU:HB3	1:A:197:LEU:HD23	1.79	0.63
1:A:59:ARG:NH2	1:A:302:GLU:OE2	2.31	0.63
2:C:207:LYS:HD2	2:C:229:ALA:HA	1.80	0.63
1:A:45:THR:O	1:A:49:GLU:N	2.32	0.62
2:C:40:SER:OG	2:C:158:ASP:OD2	2.15	0.62
2:C:222:THR:HG22	2:C:224:PRO:HD2	1.81	0.62
2:D:76:ALA:HB1	2:D:140:VAL:HG22	1.82	0.62
1:B:59:ARG:NH2	1:B:302:GLU:OE2	2.32	0.61
2:C:140:VAL:HG13	2:C:155:LEU:HD11	1.81	0.61
3:F:160:LEU:HD21	3:F:180:LEU:HD13	1.83	0.61
2:D:205:LEU:HD22	2:D:212:LEU:HB2	1.83	0.61
2:D:222:THR:HG22	2:D:224:PRO:HD2	1.83	0.61
2:D:140:VAL:HG13	2:D:155:LEU:HD11	1.82	0.60
1:B:167:SER:HB3	1:B:171:ASP:HB2	1.82	0.60
1:B:177:TYR:HE2	3:F:67:GLN:HA	1.66	0.60
2:C:76:ALA:HB1	2:C:140:VAL:HG22	1.83	0.60
1:B:33:ALA:H	1:B:56:ARG:NH1	2.01	0.59
2:C:205:LEU:HD22	2:C:212:LEU:HB2	1.85	0.58
3:F:209:PRO:HG2	3:F:232:LEU:HD21	1.85	0.58
3:F:49:SER:HB3	3:F:65:THR:HG22	1.86	0.58
1:B:45:THR:O	1:B:49:GLU:N	2.31	0.58
2:D:141:VAL:HG21	2:D:175:ILE:HD11	1.84	0.58
1:A:324:ALA:HB3	1:B:144:ARG:HH12	1.69	0.58
2:C:141:VAL:HG21	2:C:175:ILE:HD11	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:ARG:HG3	1:A:177:TYR:HE1	1.70	0.57
1:B:185:GLY:HA3	3:F:92:ARG:HB3	1.86	0.57
3:F:172:LYS:HD2	3:F:190:LYS:O	2.03	0.57
2:D:36:GLY:H	4:D:301:ANP:HNB1	1.53	0.57
1:A:33:ALA:H	1:A:56:ARG:NH1	2.02	0.57
2:D:2:SER:HB2	2:D:25:ALA:HB2	1.86	0.56
2:D:38:GLY:HA2	4:D:301:ANP:H5'1	1.87	0.56
1:A:49:GLU:HA	1:A:53:TRP:HD1	1.70	0.56
3:F:165:ASN:HB3	3:F:166:PRO:HD3	1.88	0.56
3:F:53:TYR:CG	3:F:250:ARG:HD2	2.42	0.55
3:F:243:TRP:HE1	3:F:250:ARG:HH21	1.53	0.55
1:B:160:MET:O	1:B:164:ILE:HG13	2.06	0.55
2:C:5:MET:HE3	2:C:46:MET:HB3	1.89	0.55
2:C:204:TRP:CD1	2:C:214:SER:HB2	2.42	0.55
1:A:173:ARG:HG3	1:A:177:TYR:CE1	2.42	0.54
1:B:49:GLU:HA	1:B:53:TRP:HD1	1.71	0.54
1:B:218:GLU:HG2	1:B:228:LEU:HD11	1.89	0.54
1:A:165:TYR:OH	1:B:304:PRO:HG3	2.07	0.54
2:D:171:ALA:O	2:D:175:ILE:HG22	2.08	0.54
2:C:2:SER:HB2	2:C:25:ALA:HB2	1.90	0.54
1:B:200:VAL:HG21	1:B:241:TRP:CD1	2.43	0.54
3:F:162:PHE:HB2	3:F:168:PHE:CD2	2.43	0.54
1:A:255:PHE:CZ	1:A:313:GLY:HA3	2.43	0.53
1:A:200:VAL:HG21	1:A:241:TRP:CD1	2.44	0.53
3:F:221:ILE:N	3:F:222:PRO:HD2	2.23	0.53
2:D:94:LEU:HB3	2:D:107:LEU:HD21	1.91	0.53
2:C:38:GLY:HA2	4:C:301:ANP:H5'1	1.91	0.53
1:B:64:LEU:HD21	1:B:288:LEU:HD12	1.91	0.53
1:B:77:MET:SD	1:B:88:PRO:HA	2.48	0.53
2:C:94:LEU:HB3	2:C:107:LEU:HD21	1.91	0.53
1:A:45:THR:HG22	1:A:47:ARG:H	1.73	0.52
1:B:35:GLU:OE1	1:B:56:ARG:NH2	2.42	0.52
1:A:218:GLU:HG2	1:A:228:LEU:HD11	1.91	0.52
3:F:214:ILE:HG21	3:F:224:ILE:HG13	1.91	0.52
1:B:45:THR:HG22	1:B:47:ARG:H	1.75	0.52
3:F:243:TRP:HE1	3:F:250:ARG:NH2	2.08	0.52
1:A:189:ARG:NH1	3:F:226:GLN:O	2.42	0.52
1:B:177:TYR:CE2	3:F:67:GLN:HA	2.44	0.52
2:C:171:ALA:O	2:C:175:ILE:HG22	2.09	0.52
2:D:204:TRP:CD1	2:D:214:SER:HB2	2.45	0.52
1:A:77:MET:SD	1:A:88:PRO:HA	2.50	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:112:GLY:HA2	2:C:117:ASP:OD1	2.09	0.52
1:A:50:LEU:O	1:A:54:GLN:HB2	2.09	0.52
1:B:255:PHE:CZ	1:B:313:GLY:HA3	2.45	0.52
1:A:187:ASP:OD1	1:A:188:TRP:N	2.38	0.51
3:F:260:ASN:O	3:F:263:SER:OG	2.29	0.51
2:C:162:CYS:O	2:C:163:SER:OG	2.26	0.51
2:D:239:ILE:HG22	2:D:240:GLU:HG3	1.93	0.51
1:A:64:LEU:HD21	1:A:288:LEU:HD12	1.93	0.51
1:B:50:LEU:O	1:B:54:GLN:HB2	2.10	0.51
2:D:5:MET:HE3	2:D:46:MET:HB3	1.91	0.51
3:F:162:PHE:HB2	3:F:168:PHE:CE2	2.45	0.51
2:C:243:ARG:H	2:C:243:ARG:HD2	1.74	0.51
1:B:173:ARG:HG3	1:B:177:TYR:CE1	2.46	0.50
2:C:134:ARG:CD	2:C:168:GLN:HG3	2.41	0.50
2:D:243:ARG:H	2:D:243:ARG:HD2	1.75	0.50
2:D:112:GLY:HA2	2:D:117:ASP:OD1	2.10	0.50
2:C:239:ILE:HG22	2:C:240:GLU:HG3	1.92	0.50
1:A:35:GLU:OE1	1:A:56:ARG:NH2	2.44	0.50
2:C:35:ASN:HA	4:C:301:ANP:HNB1	1.76	0.50
1:B:158:ALA:O	1:B:162:TRP:HB2	2.12	0.50
3:F:86:ARG:HD3	3:F:91:GLU:OE2	2.12	0.50
1:A:112:LEU:HD23	1:A:113:PRO:HD3	1.94	0.50
3:F:172:LYS:H	3:F:172:LYS:HD3	1.76	0.50
2:C:166:VAL:HG11	2:D:234:PHE:HZ	1.77	0.49
3:F:53:TYR:CD2	3:F:250:ARG:HD2	2.47	0.49
1:A:165:TYR:CD1	3:F:67:GLN:HB3	2.47	0.49
3:F:155:LYS:HB3	3:F:186:GLU:HG2	1.94	0.49
1:B:101:LEU:HD21	1:B:117:LEU:HB3	1.94	0.49
1:A:170:VAL:HG13	3:F:162:PHE:HB3	1.94	0.49
2:C:234:PHE:HZ	2:D:166:VAL:HG11	1.78	0.49
1:B:72:ILE:HD11	1:B:280:ALA:HB2	1.94	0.49
1:B:173:ARG:HG3	1:B:177:TYR:HE1	1.77	0.49
2:D:5:MET:HE2	2:D:154:LEU:HD21	1.95	0.49
1:A:97:ALA:HB1	1:A:121:ALA:HB1	1.95	0.49
2:C:197:LEU:HD12	2:C:197:LEU:HA	1.67	0.49
1:B:92:GLY:O	1:B:96:GLY:N	2.46	0.48
2:D:35:ASN:HA	4:D:301:ANP:HNB1	1.78	0.48
1:B:213:MET:HE2	1:B:226:LEU:HB2	1.95	0.48
2:C:71:LEU:HD13	2:C:75:ARG:HH21	1.78	0.48
2:C:78:LEU:HD23	2:C:79:SER:O	2.13	0.48
1:A:213:MET:HE3	1:A:224:LEU:HB3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:82:GLN:NE2	2:D:136:ARG:HE	2.11	0.48
3:F:159:PHE:CE2	3:F:161:GLN:HB2	2.49	0.48
2:D:192:ASP:OD1	2:D:195:HIS:N	2.41	0.48
1:B:97:ALA:HB1	1:B:121:ALA:HB1	1.95	0.48
1:A:158:ALA:O	1:A:162:TRP:HB2	2.14	0.47
1:B:38:ILE:HG23	1:B:42:ASP:HB3	1.96	0.47
1:B:157:SER:O	1:B:161:THR:HB	2.13	0.47
2:C:10:VAL:O	2:C:18:PRO:HA	2.14	0.47
1:A:173:ARG:CZ	3:F:196:TRP:HB2	2.44	0.47
1:A:72:ILE:HD11	1:A:280:ALA:HB2	1.95	0.47
1:A:303:LEU:HD12	1:A:303:LEU:HA	1.77	0.47
2:C:82:GLN:NE2	2:C:136:ARG:HE	2.12	0.47
1:B:187:ASP:OD1	1:B:188:TRP:N	2.45	0.47
1:B:213:MET:HE3	1:B:224:LEU:HB3	1.97	0.47
2:D:80:GLN:HB3	2:D:158:ASP:OD2	2.14	0.47
1:A:173:ARG:HD2	1:A:173:ARG:HA	1.63	0.47
3:F:82:VAL:HB	3:F:104:VAL:HG22	1.97	0.47
1:A:95:ASN:HB2	1:A:156:SER:HB3	1.97	0.47
2:C:82:GLN:HE21	2:C:136:ARG:HH21	1.62	0.47
1:A:92:GLY:HA3	1:A:153:GLY:HA2	1.97	0.47
1:A:157:SER:O	1:A:161:THR:HB	2.15	0.47
1:B:112:LEU:HD23	1:B:113:PRO:HD3	1.97	0.47
2:C:5:MET:HE2	2:C:154:LEU:HD21	1.97	0.47
1:A:38:ILE:HG23	1:A:42:ASP:HB3	1.96	0.46
1:A:160:MET:O	1:A:164:ILE:HG13	2.16	0.46
1:A:101:LEU:HD21	1:A:117:LEU:HB3	1.96	0.46
2:D:82:GLN:HE21	2:D:136:ARG:HH21	1.63	0.46
2:D:78:LEU:HD23	2:D:79:SER:O	2.15	0.46
2:C:80:GLN:HB3	2:C:158:ASP:OD2	2.15	0.46
3:F:200:SER:O	3:F:204:VAL:HG23	2.16	0.46
3:F:166:PRO:HA	3:F:167:PRO:HD3	1.82	0.46
1:A:213:MET:HE2	1:A:226:LEU:HB2	1.98	0.46
3:F:94:VAL:CG1	3:F:104:VAL:HG11	2.45	0.46
3:F:32:ALA:HB2	3:F:246:ARG:HB3	1.97	0.45
1:B:165:TYR:HE1	3:F:194:VAL:HG21	1.82	0.45
2:C:217:ARG:HH21	2:C:243:ARG:NH1	2.14	0.45
1:B:165:TYR:CE1	3:F:194:VAL:HG21	2.51	0.45
2:C:159:GLN:H	2:C:189:SER:HB2	1.81	0.45
2:D:71:LEU:HD13	2:D:75:ARG:HH21	1.81	0.45
2:C:118:ASP:OD1	2:C:118:ASP:N	2.38	0.45
1:A:165:TYR:HD1	3:F:67:GLN:HE21	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:10:VAL:O	2:D:18:PRO:HA	2.17	0.45
1:A:92:GLY:O	1:A:96:GLY:N	2.47	0.45
1:B:92:GLY:HA3	1:B:153:GLY:HA2	1.98	0.45
1:B:173:ARG:HD2	1:B:173:ARG:HA	1.64	0.45
1:B:287:LEU:HD23	1:B:287:LEU:HA	1.71	0.45
3:F:28:THR:HG22	3:F:83:ILE:HD12	1.99	0.45
3:F:160:LEU:HD23	3:F:213:VAL:HB	1.98	0.45
2:C:192:ASP:OD1	2:C:195:HIS:N	2.42	0.45
3:F:69:MET:HG2	3:F:93:GLN:CD	2.42	0.45
2:C:16:LEU:HD22	2:C:17:GLY:O	2.17	0.44
1:A:220:SER:OG	2:C:82:GLN:OE1	2.18	0.44
2:C:43:LEU:HD12	2:C:43:LEU:HA	1.82	0.44
2:C:36:GLY:H	4:C:301:ANP:HNB1	1.66	0.44
2:D:159:GLN:H	2:D:189:SER:HB2	1.82	0.44
2:D:33:GLY:O	2:D:39:LYS:NZ	2.49	0.44
1:B:95:ASN:HB2	1:B:156:SER:HB3	1.99	0.44
1:A:287:LEU:HD23	1:A:287:LEU:HA	1.69	0.43
1:B:54:GLN:HA	1:B:188:TRP:NE1	2.31	0.43
1:B:78:GLN:OE1	1:B:265:ARG:NH2	2.48	0.43
1:A:302:GLU:OE1	3:F:200:SER:HB2	2.19	0.43
1:A:78:GLN:OE1	1:A:265:ARG:NH2	2.48	0.43
3:F:232:LEU:HG	3:F:234:ILE:HD13	1.99	0.43
2:C:43:LEU:HD23	2:C:158:ASP:HB2	2.01	0.43
2:C:101:LYS:H	2:C:101:LYS:HG2	1.60	0.43
2:D:82:GLN:HE21	2:D:136:ARG:HE	1.67	0.43
4:C:301:ANP:O2'	2:D:122:ARG:NH1	2.46	0.43
1:A:152:LEU:HD23	1:A:152:LEU:HA	1.80	0.42
2:C:19:LEU:HD22	2:C:211:MET:HB2	1.99	0.42
1:B:228:LEU:HA	1:B:228:LEU:HD23	1.72	0.42
2:D:197:LEU:HD12	2:D:197:LEU:HA	1.68	0.42
3:F:217:GLY:C	3:F:219:ASP:H	2.27	0.42
2:C:105:GLU:CD	2:C:105:GLU:H	2.27	0.42
2:D:19:LEU:HD22	2:D:211:MET:HB2	2.00	0.42
3:F:157:ARG:HH21	3:F:188:ILE:HA	1.83	0.42
1:B:95:ASN:CB	1:B:156:SER:HB3	2.49	0.42
1:A:228:LEU:HD23	1:A:228:LEU:HA	1.69	0.42
1:B:298:LEU:HD12	1:B:298:LEU:H	1.84	0.42
4:C:301:ANP:N6	2:D:126:GLN:HG2	2.35	0.42
1:B:11:GLN:HG2	1:B:14:ARG:HH21	1.85	0.42
1:B:55:ILE:HD11	3:F:100:LEU:HD21	2.02	0.42
1:A:29:LEU:HD22	1:A:29:LEU:HA	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298:LEU:HD21	1:B:303:LEU:HD13	2.02	0.42
2:C:172:LEU:HA	2:C:175:ILE:CG2	2.50	0.42
3:F:194:VAL:HB	3:F:197:PRO:HB3	2.01	0.42
1:A:95:ASN:CB	1:A:156:SER:HB3	2.48	0.42
1:A:173:ARG:O	1:A:177:TYR:HD1	2.02	0.42
3:F:47:VAL:HG23	3:F:62:GLN:HG2	2.02	0.42
1:A:112:LEU:HD12	3:F:222:PRO:HB2	2.01	0.41
1:A:224:LEU:HD11	2:C:97:HIS:HA	2.02	0.41
2:C:114:LEU:HD12	2:C:138:ALA:HB2	2.02	0.41
3:F:97:LEU:HD22	3:F:102:ILE:HD12	2.01	0.41
1:A:165:TYR:HE1	1:B:304:PRO:CD	2.32	0.41
1:A:54:GLN:HA	1:A:188:TRP:NE1	2.31	0.41
4:C:301:ANP:HN62	2:D:126:GLN:HG2	1.85	0.41
2:D:115:ALA:HB3	2:D:134:ARG:NH1	2.36	0.41
3:F:29:LEU:HD12	3:F:84:ALA:HB2	2.02	0.41
3:F:209:PRO:CG	3:F:232:LEU:HD21	2.48	0.41
2:D:165:ASP:OD1	2:D:168:GLN:HG2	2.21	0.41
2:D:217:ARG:HH21	2:D:243:ARG:NH1	2.18	0.41
1:A:11:GLN:HG2	1:A:14:ARG:HH21	1.85	0.41
1:A:172:LEU:HD22	1:A:172:LEU:HA	1.95	0.41
1:B:173:ARG:O	1:B:177:TYR:HD1	2.03	0.41
1:B:129:THR:O	1:B:133:LEU:HG	2.21	0.41
2:D:114:LEU:HD12	2:D:138:ALA:HB2	2.03	0.41
2:D:159:GLN:HG3	2:D:189:SER:HB2	2.03	0.41
3:F:122:LEU:HD23	3:F:122:LEU:HA	1.91	0.41
2:D:105:GLU:CD	2:D:105:GLU:H	2.28	0.41
1:B:321:LEU:HD23	1:B:321:LEU:HA	1.94	0.40
3:F:94:VAL:HG11	3:F:104:VAL:HG11	2.03	0.40
1:A:30:SER:HB2	1:A:295:ARG:HH22	1.86	0.40
1:A:125:ALA:O	1:A:129:THR:HG23	2.21	0.40
1:B:226:LEU:HA	1:B:227:PRO:HD2	1.87	0.40
2:C:159:GLN:HG3	2:C:189:SER:HB2	2.03	0.40
2:D:35:ASN:HA	4:D:301:ANP:N3B	2.36	0.40
1:A:161:THR:HG23	1:B:180:MET:HE3	2.02	0.40
1:B:113:PRO:HG2	1:B:116:ALA:HB3	2.03	0.40
2:C:82:GLN:HE21	2:C:136:ARG:HE	1.68	0.40
2:C:122:ARG:NH1	4:D:301:ANP:O2'	2.48	0.40
1:B:57:LEU:HA	1:B:60:THR:HG22	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	322/349 (92%)	303 (94%)	19 (6%)	0	100	100
1	B	322/349 (92%)	303 (94%)	19 (6%)	0	100	100
2	C	246/249 (99%)	238 (97%)	7 (3%)	1 (0%)	30	62
2	D	246/249 (99%)	238 (97%)	8 (3%)	0	100	100
3	F	243/255 (95%)	229 (94%)	11 (4%)	3 (1%)	10	41
All	All	1379/1451 (95%)	1311 (95%)	64 (5%)	4 (0%)	36	67

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	230	GLU
3	F	233	LYS
2	C	166	VAL
3	F	166	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	249/270 (92%)	227 (91%)	22 (9%)	9	33
1	B	249/270 (92%)	227 (91%)	22 (9%)	9	33
2	C	200/201 (100%)	177 (88%)	23 (12%)	5	25
2	D	200/201 (100%)	178 (89%)	22 (11%)	6	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	F	205/214 (96%)	200 (98%)	5 (2%)	43 65
All	All	1103/1156 (95%)	1009 (92%)	94 (8%)	10 35

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	20	SER
1	A	29	LEU
1	A	112	LEU
1	A	114	ASN
1	A	115	TRP
1	A	142	THR
1	A	156	SER
1	A	161	THR
1	A	171	ASP
1	A	172	LEU
1	A	173	ARG
1	A	195	LEU
1	A	211	MET
1	A	218	GLU
1	A	220	SER
1	A	224	LEU
1	A	274	VAL
1	A	287	LEU
1	A	288	LEU
1	A	305	ILE
1	A	308	VAL
1	B	8	GLN
1	B	20	SER
1	B	29	LEU
1	B	112	LEU
1	B	114	ASN
1	B	115	TRP
1	B	142	THR
1	B	156	SER
1	B	161	THR
1	B	171	ASP
1	B	172	LEU
1	B	173	ARG
1	B	195	LEU
1	B	211	MET

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Mol	Chain	Res	Type
1	B	218	GLU
1	B	220	SER
1	B	224	LEU
1	B	274	VAL
1	B	287	LEU
1	B	288	LEU
1	B	305	ILE
1	B	308	VAL
2	C	15	ARG
2	C	16	LEU
2	C	32	VAL
2	C	50	THR
2	C	56	ILE
2	C	71	LEU
2	C	96	LEU
2	C	100	ASP
2	C	104	THR
2	C	105	GLU
2	C	106	LEU
2	C	108	ASN
2	C	118	ASP
2	C	128	SER
2	C	140	VAL
2	C	162	CYS
2	C	168	GLN
2	C	175	ILE
2	C	190	SER
2	C	191	HIS
2	C	197	LEU
2	C	217	ARG
2	C	243	ARG
2	D	15	ARG
2	D	16	LEU
2	D	32	VAL
2	D	50	THR
2	D	56	ILE
2	D	71	LEU
2	D	96	LEU
2	D	100	ASP
2	D	104	THR
2	D	105	GLU
2	D	106	LEU

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Mol	Chain	Res	Type
2	D	108	ASN
2	D	118	ASP
2	D	128	SER
2	D	140	VAL
2	D	162	CYS
2	D	175	ILE
2	D	190	SER
2	D	191	HIS
2	D	197	LEU
2	D	217	ARG
2	D	243	ARG
3	F	67	GLN
3	F	83	ILE
3	F	164	ILE
3	F	205	LEU
3	F	228	TRP

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

Mol	Chain	Res	Type
2	C	182	GLN
2	D	81	GLN
2	D	182	GLN
3	F	128	GLN
3	F	198	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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ANP	D	301	5	33,33,33	2.05	12 (36%)	45,52,52	2.24	12 (26%)
4	ANP	C	301	5	33,33,33	2.13	13 (39%)	45,52,52	2.29	13 (28%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ANP	D	301	5	-	9/18/38/38	0/3/3/3
4	ANP	C	301	5	-	9/18/38/38	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	301	ANP	C5-C4	4.77	1.47	1.39
4	C	301	ANP	C5-C4	4.75	1.47	1.39
4	C	301	ANP	PG-N3B	4.51	1.75	1.63
4	D	301	ANP	PB-N3B	4.47	1.75	1.63
4	C	301	ANP	PB-N3B	4.46	1.75	1.63
4	D	301	ANP	PG-N3B	4.43	1.74	1.63
4	C	301	ANP	PG-O1G	3.68	1.51	1.46
4	C	301	ANP	PB-O1B	3.64	1.51	1.46
4	D	301	ANP	PG-O1G	3.44	1.51	1.46
4	D	301	ANP	PB-O1B	3.18	1.51	1.46
4	D	301	ANP	PB-O2B	-2.84	1.49	1.56
4	C	301	ANP	PB-O2B	-2.81	1.49	1.56
4	C	301	ANP	C4-N9	-2.66	1.32	1.37
4	D	301	ANP	C8-N7	2.65	1.36	1.31
4	D	301	ANP	C4-N9	-2.64	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	301	ANP	PG-O3G	-2.61	1.49	1.56
4	C	301	ANP	C5-C6	2.58	1.48	1.41
4	D	301	ANP	C5-C6	2.57	1.48	1.41
4	C	301	ANP	PG-O3G	-2.50	1.50	1.56
4	C	301	ANP	C8-N7	2.49	1.36	1.31
4	C	301	ANP	C5-N7	-2.44	1.34	1.39
4	D	301	ANP	PG-O2G	-2.34	1.50	1.56
4	D	301	ANP	C5-N7	-2.32	1.34	1.39
4	C	301	ANP	PB-O3A	2.32	1.62	1.59
4	C	301	ANP	PG-O2G	-2.25	1.50	1.56

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	301	ANP	O1G-PG-N3B	-9.10	98.37	111.77
4	D	301	ANP	O1G-PG-N3B	-9.08	98.40	111.77
4	D	301	ANP	C5-C4-N3	-4.86	120.03	126.72
4	C	301	ANP	C5-C4-N3	-4.58	120.41	126.72
4	C	301	ANP	O1B-PB-N3B	-4.45	105.22	111.77
4	D	301	ANP	O2B-PB-O1B	3.67	117.73	109.87
4	D	301	ANP	N3-C4-N9	3.53	133.17	127.17
4	C	301	ANP	N3-C4-N9	3.48	133.08	127.17
4	D	301	ANP	O1B-PB-N3B	-3.45	106.70	111.77
4	C	301	ANP	N3-C2-N1	-3.33	123.54	128.58
4	D	301	ANP	C2-N3-C4	3.31	119.93	111.83
4	C	301	ANP	C2-N3-C4	3.28	119.85	111.83
4	C	301	ANP	O2B-PB-O1B	3.09	116.50	109.87
4	D	301	ANP	N3-C2-N1	-3.02	124.01	128.58
4	C	301	ANP	C4-C5-N7	-2.97	107.19	110.58
4	D	301	ANP	C4-C5-N7	-2.96	107.19	110.58
4	C	301	ANP	O3G-PG-O1G	2.68	120.17	113.45
4	D	301	ANP	O3G-PG-O1G	2.50	119.71	113.45
4	C	301	ANP	C4-N9-C8	2.46	108.33	105.74
4	C	301	ANP	C2-N1-C6	2.29	122.49	118.73
4	C	301	ANP	C6-C5-N7	2.16	136.25	132.09
4	D	301	ANP	C4-N9-C8	2.15	108.00	105.74
4	C	301	ANP	C2'-C3'-C4'	2.04	106.54	102.61
4	D	301	ANP	C2-N1-C6	2.03	122.06	118.73
4	D	301	ANP	C6-C5-N7	2.03	136.00	132.09

There are no chirality outliers.

All (18) torsion outliers are listed below:

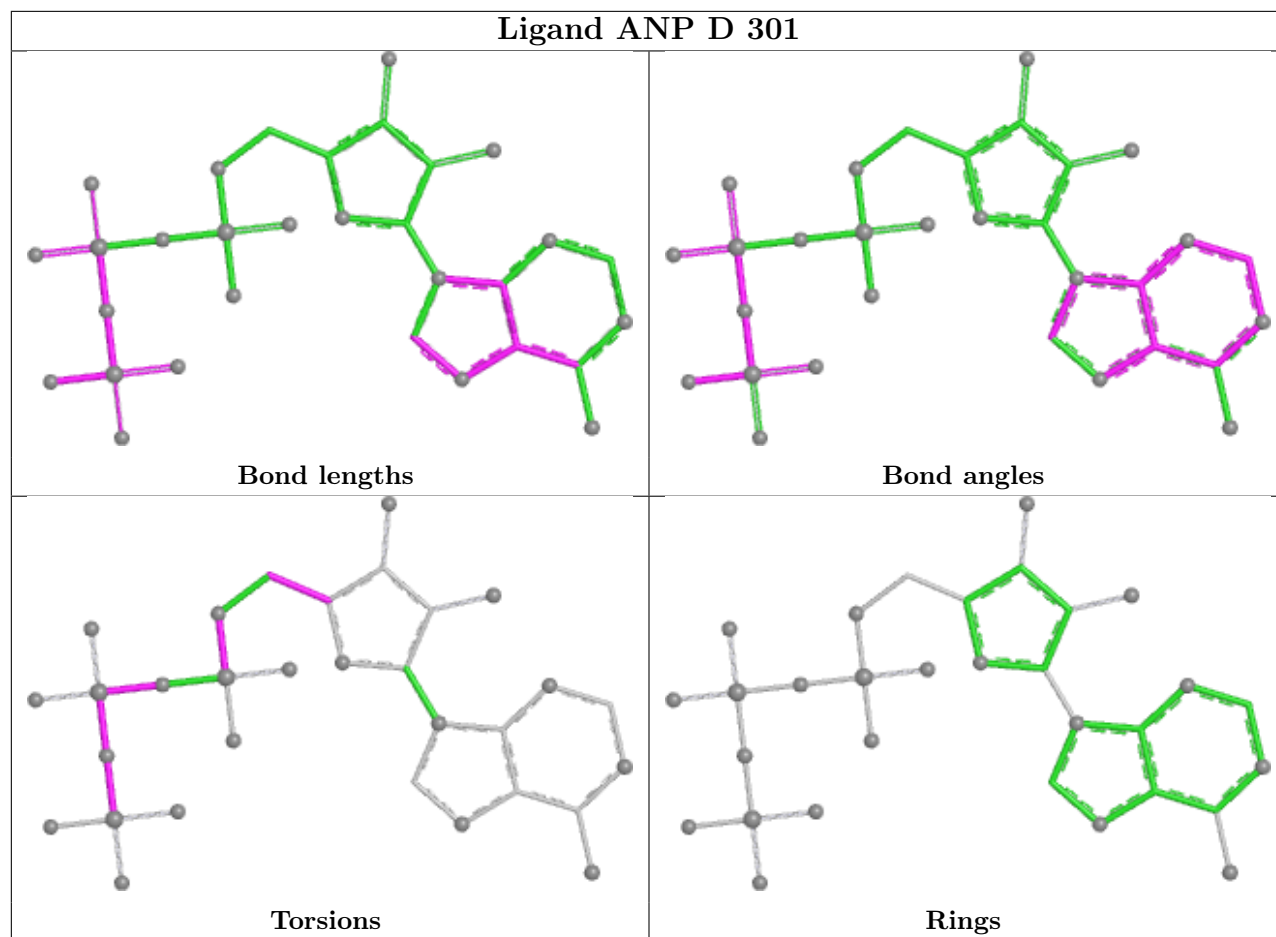
Mol	Chain	Res	Type	Atoms
4	C	301	ANP	PB-N3B-PG-O1G
4	C	301	ANP	PG-N3B-PB-O1B
4	C	301	ANP	PA-O3A-PB-O2B
4	C	301	ANP	C5'-O5'-PA-O1A
4	D	301	ANP	PB-N3B-PG-O1G
4	D	301	ANP	PG-N3B-PB-O1B
4	D	301	ANP	PA-O3A-PB-O2B
4	D	301	ANP	C5'-O5'-PA-O1A
4	C	301	ANP	O4'-C4'-C5'-O5'
4	D	301	ANP	O4'-C4'-C5'-O5'
4	C	301	ANP	C3'-C4'-C5'-O5'
4	D	301	ANP	C3'-C4'-C5'-O5'
4	C	301	ANP	C5'-O5'-PA-O2A
4	C	301	ANP	C5'-O5'-PA-O3A
4	D	301	ANP	C5'-O5'-PA-O2A
4	D	301	ANP	C5'-O5'-PA-O3A
4	C	301	ANP	PA-O3A-PB-O1B
4	D	301	ANP	PA-O3A-PB-O1B

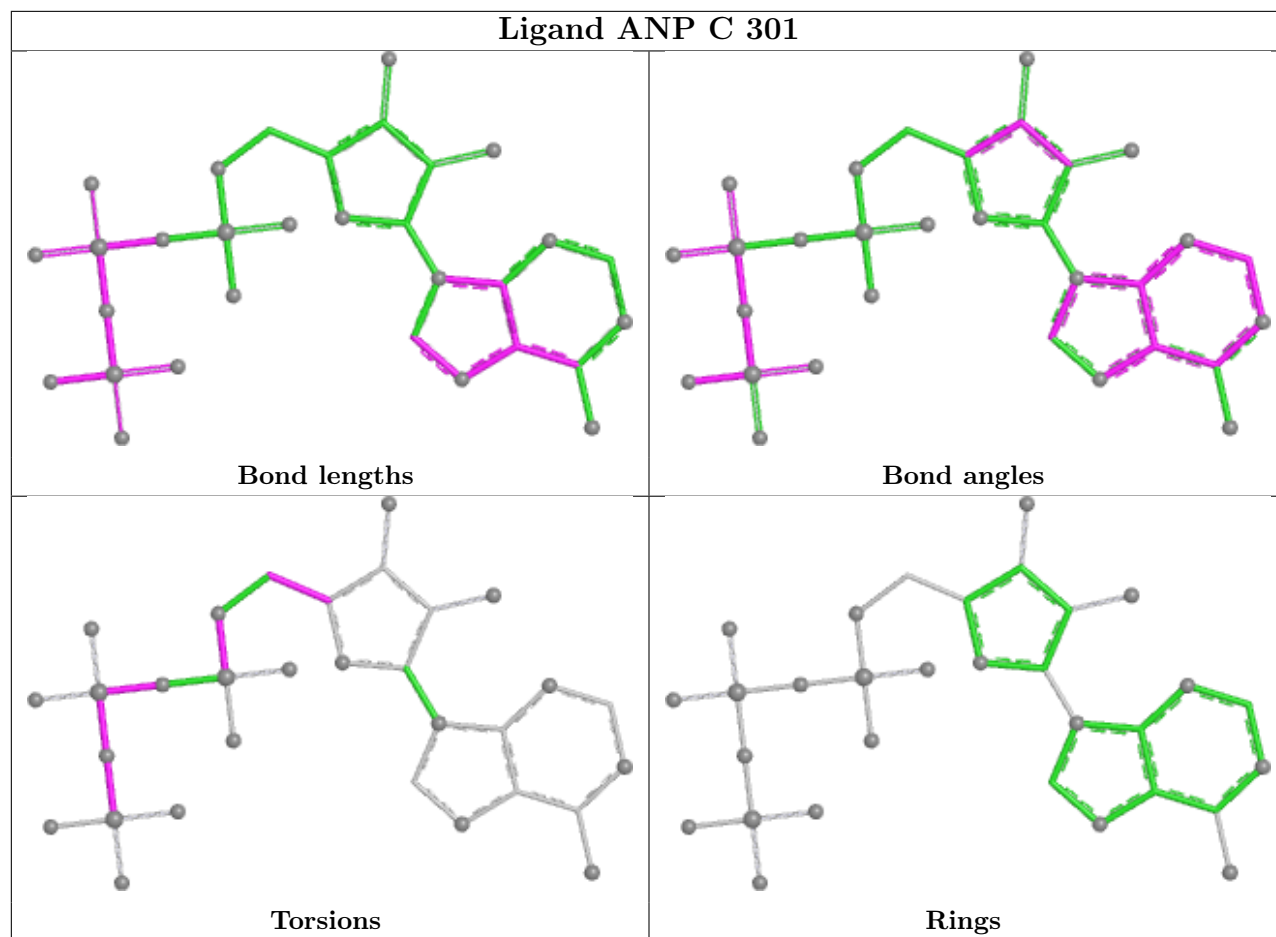
There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	301	ANP	5	0
4	C	301	ANP	6	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	324/349 (92%)	-0.13	6 (1%) 66 42	85, 144, 237, 373	0
1	B	324/349 (92%)	-0.09	5 (1%) 72 48	103, 151, 239, 405	0
2	C	248/249 (99%)	-0.34	2 (0%) 82 61	68, 113, 195, 311	0
2	D	248/249 (99%)	-0.35	1 (0%) 88 73	79, 134, 200, 262	0
3	F	245/255 (96%)	0.40	18 (7%) 21 15	173, 255, 350, 525	0
All	All	1389/1451 (95%)	-0.11	32 (2%) 61 38	68, 148, 286, 525	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	164	ILE	8.6
3	F	205	LEU	5.4
3	F	94	VAL	4.8
3	F	163	GLY	3.9
3	F	166	PRO	3.9
3	F	204	VAL	3.7
1	B	190	GLN	3.7
1	A	93	VAL	3.5
1	B	176	MET	3.5
3	F	104	VAL	3.4
1	A	178	TRP	3.3
3	F	165	ASN	3.1
3	F	112	ILE	3.0
1	A	51	PHE	3.0
3	F	198	GLN	2.9
3	F	247	ALA	2.8
3	F	139	LEU	2.5
1	A	2	LEU	2.5
3	F	92	ARG	2.5
3	F	159	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	50	LEU	2.3
2	D	190	SER	2.3
3	F	232	LEU	2.3
3	F	33	ASN	2.3
1	A	176	MET	2.2
2	C	190	SER	2.2
2	C	192	ASP	2.2
1	B	88	PRO	2.1
1	B	236	VAL	2.1
3	F	249	PRO	2.1
3	F	189	PHE	2.0
1	A	85	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

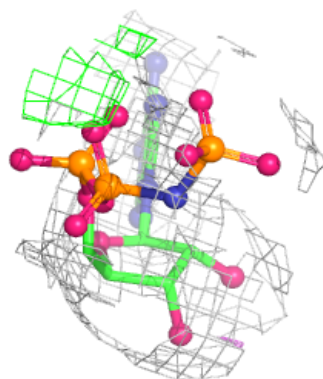
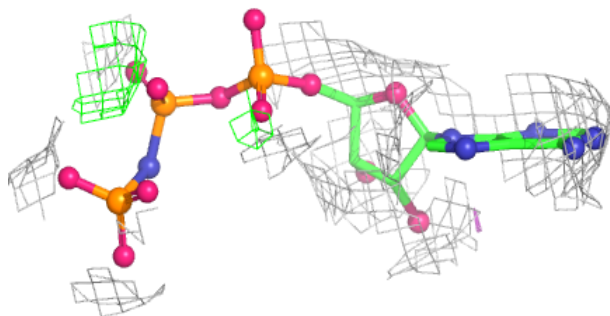
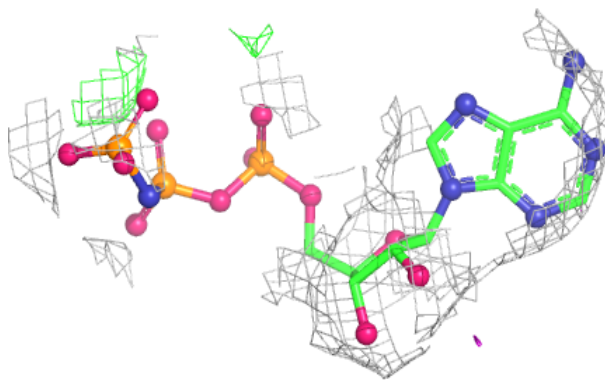
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ANP	D	301	31/31	0.94	0.07	101,123,145,146	0
4	ANP	C	301	31/31	0.95	0.08	101,120,141,142	0
5	MG	D	302	1/1	0.95	0.07	57,57,57,57	0
5	MG	C	302	1/1	0.96	0.06	57,57,57,57	0

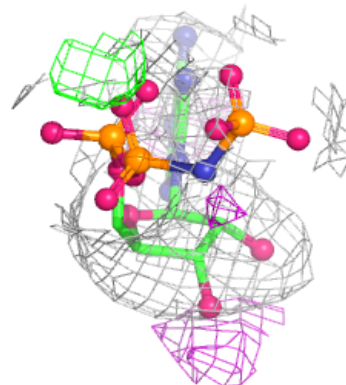
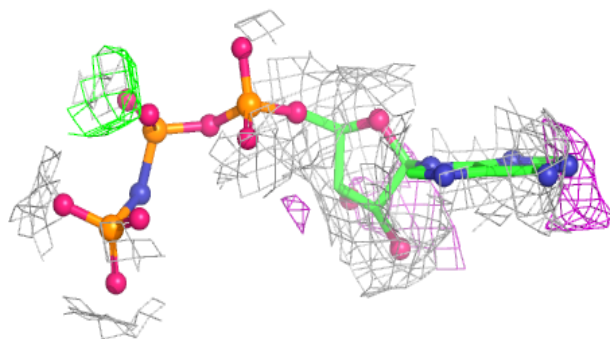
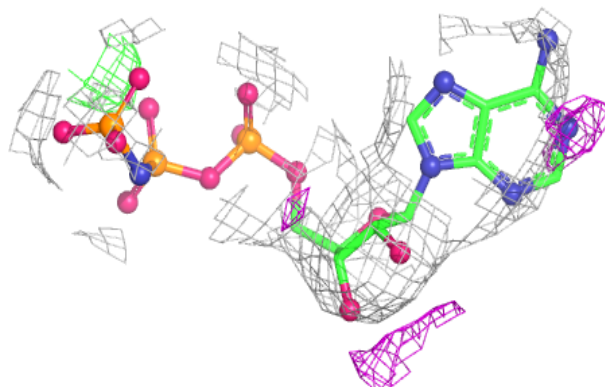
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP D 301:

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

**Electron density around ANP C 301:**

$2mF_o-DF_c$ (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.