



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:11 AM UTC

PDB ID : 2OJW / pdb_00002ojw
Title : Crystal structure of human glutamine synthetase in complex with ADP and phosphate
Authors : Karlberg, T.; Uppenberg, J.; Arrowsmith, C.; Berglund, H.; Busam, R.D.; Collins, R.; Edwards, A.; Flodin, S.; Flores, A.; Graslund, S.; Hallberg, B.M.; Hammarstrom, M.; Hogbom, M.; Johansson, I.; Kotenyova, T.; Moche, M.; Nilsson, M.E.; Nordlund, P.; Nyman, T.; Ogg, D.; Persson, C.; Sagemark, J.; Stenmark, P.; Sundstrom, M.; Thorsell, A.G.; Van Den Berg, S.; Wallden, K.; Weigelt, J.; Holmberg-Schiavone, L.; Structural Genomics Consortium (SGC)
Deposited on : 2007-01-15
Resolution : 2.05 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)

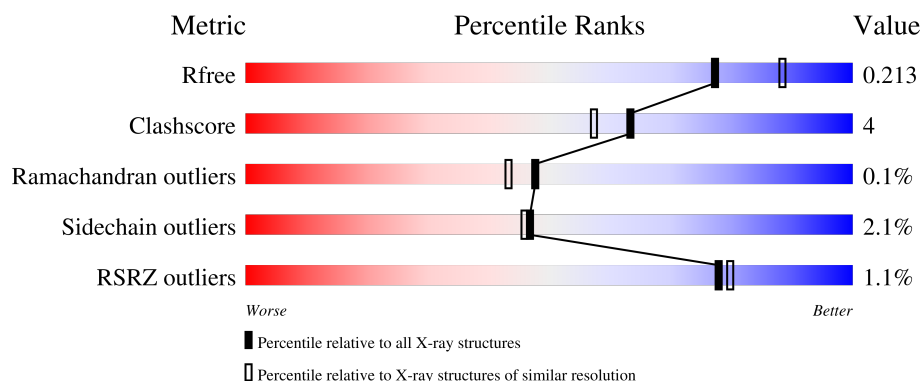
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	384	
1	B	384	

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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

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Mol	Chain	Length	Quality of chain
1	C	384	% 85%9%5%
1	D	384	% 84%10%6%
1	E	384	83%8%8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15882 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 Glutamine synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	1	0
			2855	1795	508	528	24			
1	B	362	Total	C	N	O	S	0	1	0
			2870	1802	513	532	23			
1	C	364	Total	C	N	O	S	0	0	0
			2874	1804	512	535	23			
1	D	362	Total	C	N	O	S	0	0	0
			2863	1798	510	532	23			
1	E	354	Total	C	N	O	S	0	1	0
			2809	1763	504	520	22			

There are 115 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	expression tag	UNP P15104
A	-17	HIS	-	expression tag	UNP P15104
A	-16	HIS	-	expression tag	UNP P15104
A	-15	HIS	-	expression tag	UNP P15104
A	-14	HIS	-	expression tag	UNP P15104
A	-13	HIS	-	expression tag	UNP P15104
A	-12	HIS	-	expression tag	UNP P15104
A	-11	SER	-	expression tag	UNP P15104
A	-10	SER	-	expression tag	UNP P15104
A	-9	GLY	-	expression tag	UNP P15104
A	-8	VAL	-	expression tag	UNP P15104
A	-7	ASP	-	expression tag	UNP P15104
A	-6	LEU	-	expression tag	UNP P15104
A	-5	GLY	-	expression tag	UNP P15104
A	-4	THR	-	expression tag	UNP P15104
A	-3	GLU	-	expression tag	UNP P15104
A	-2	ASN	-	expression tag	UNP P15104
A	-1	LEU	-	expression tag	UNP P15104
A	0	TYR	-	expression tag	UNP P15104

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1	PHE	-	expression tag	UNP P15104
A	2	GLN	-	expression tag	UNP P15104
A	3	SER	-	expression tag	UNP P15104
A	4	MET	-	expression tag	UNP P15104
B	-18	MET	-	expression tag	UNP P15104
B	-17	HIS	-	expression tag	UNP P15104
B	-16	HIS	-	expression tag	UNP P15104
B	-15	HIS	-	expression tag	UNP P15104
B	-14	HIS	-	expression tag	UNP P15104
B	-13	HIS	-	expression tag	UNP P15104
B	-12	HIS	-	expression tag	UNP P15104
B	-11	SER	-	expression tag	UNP P15104
B	-10	SER	-	expression tag	UNP P15104
B	-9	GLY	-	expression tag	UNP P15104
B	-8	VAL	-	expression tag	UNP P15104
B	-7	ASP	-	expression tag	UNP P15104
B	-6	LEU	-	expression tag	UNP P15104
B	-5	GLY	-	expression tag	UNP P15104
B	-4	THR	-	expression tag	UNP P15104
B	-3	GLU	-	expression tag	UNP P15104
B	-2	ASN	-	expression tag	UNP P15104
B	-1	LEU	-	expression tag	UNP P15104
B	0	TYR	-	expression tag	UNP P15104
B	1	PHE	-	expression tag	UNP P15104
B	2	GLN	-	expression tag	UNP P15104
B	3	SER	-	expression tag	UNP P15104
B	4	MET	-	expression tag	UNP P15104
C	-18	MET	-	expression tag	UNP P15104
C	-17	HIS	-	expression tag	UNP P15104
C	-16	HIS	-	expression tag	UNP P15104
C	-15	HIS	-	expression tag	UNP P15104
C	-14	HIS	-	expression tag	UNP P15104
C	-13	HIS	-	expression tag	UNP P15104
C	-12	HIS	-	expression tag	UNP P15104
C	-11	SER	-	expression tag	UNP P15104
C	-10	SER	-	expression tag	UNP P15104
C	-9	GLY	-	expression tag	UNP P15104
C	-8	VAL	-	expression tag	UNP P15104
C	-7	ASP	-	expression tag	UNP P15104
C	-6	LEU	-	expression tag	UNP P15104
C	-5	GLY	-	expression tag	UNP P15104
C	-4	THR	-	expression tag	UNP P15104

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLU	-	expression tag	UNP P15104
C	-2	ASN	-	expression tag	UNP P15104
C	-1	LEU	-	expression tag	UNP P15104
C	0	TYR	-	expression tag	UNP P15104
C	1	PHE	-	expression tag	UNP P15104
C	2	GLN	-	expression tag	UNP P15104
C	3	SER	-	expression tag	UNP P15104
C	4	MET	-	expression tag	UNP P15104
D	-18	MET	-	expression tag	UNP P15104
D	-17	HIS	-	expression tag	UNP P15104
D	-16	HIS	-	expression tag	UNP P15104
D	-15	HIS	-	expression tag	UNP P15104
D	-14	HIS	-	expression tag	UNP P15104
D	-13	HIS	-	expression tag	UNP P15104
D	-12	HIS	-	expression tag	UNP P15104
D	-11	SER	-	expression tag	UNP P15104
D	-10	SER	-	expression tag	UNP P15104
D	-9	GLY	-	expression tag	UNP P15104
D	-8	VAL	-	expression tag	UNP P15104
D	-7	ASP	-	expression tag	UNP P15104
D	-6	LEU	-	expression tag	UNP P15104
D	-5	GLY	-	expression tag	UNP P15104
D	-4	THR	-	expression tag	UNP P15104
D	-3	GLU	-	expression tag	UNP P15104
D	-2	ASN	-	expression tag	UNP P15104
D	-1	LEU	-	expression tag	UNP P15104
D	0	TYR	-	expression tag	UNP P15104
D	1	PHE	-	expression tag	UNP P15104
D	2	GLN	-	expression tag	UNP P15104
D	3	SER	-	expression tag	UNP P15104
D	4	MET	-	expression tag	UNP P15104
E	-18	MET	-	expression tag	UNP P15104
E	-17	HIS	-	expression tag	UNP P15104
E	-16	HIS	-	expression tag	UNP P15104
E	-15	HIS	-	expression tag	UNP P15104
E	-14	HIS	-	expression tag	UNP P15104
E	-13	HIS	-	expression tag	UNP P15104
E	-12	HIS	-	expression tag	UNP P15104
E	-11	SER	-	expression tag	UNP P15104
E	-10	SER	-	expression tag	UNP P15104
E	-9	GLY	-	expression tag	UNP P15104
E	-8	VAL	-	expression tag	UNP P15104

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-7	ASP	-	expression tag	UNP P15104
E	-6	LEU	-	expression tag	UNP P15104
E	-5	GLY	-	expression tag	UNP P15104
E	-4	THR	-	expression tag	UNP P15104
E	-3	GLU	-	expression tag	UNP P15104
E	-2	ASN	-	expression tag	UNP P15104
E	-1	LEU	-	expression tag	UNP P15104
E	0	TYR	-	expression tag	UNP P15104
E	1	PHE	-	expression tag	UNP P15104
E	2	GLN	-	expression tag	UNP P15104
E	3	SER	-	expression tag	UNP P15104
E	4	MET	-	expression tag	UNP P15104

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Mn 4 4	0	0
2	B	4	Total Mn 4 4	0	0
2	C	4	Total Mn 4 4	0	0
2	D	4	Total Mn 4 4	0	0
2	E	4	Total Mn 4 4	0	0

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

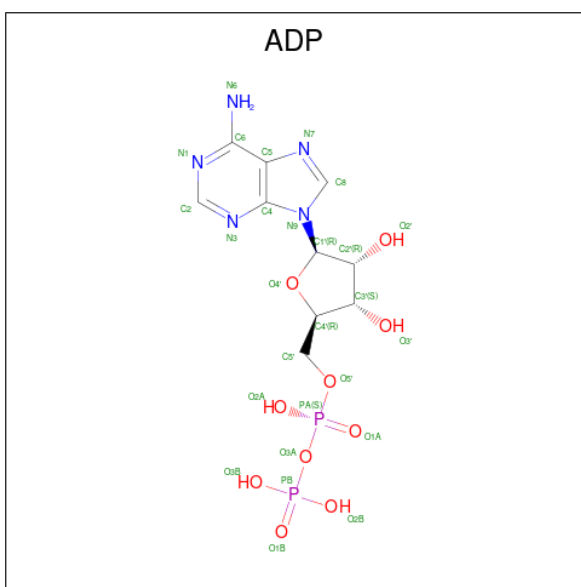


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	E	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

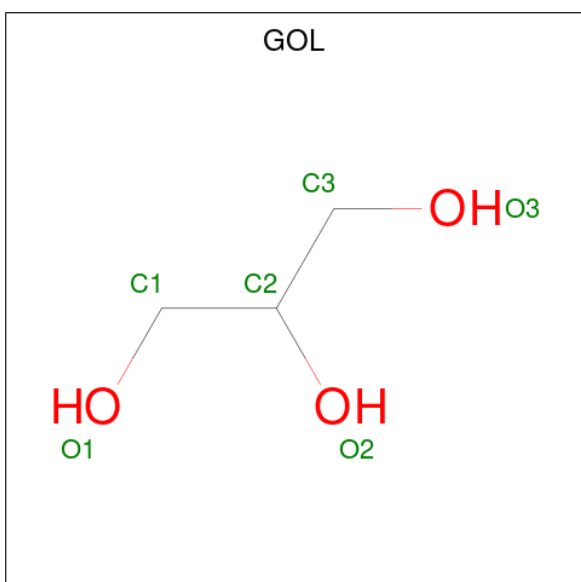
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Cl	0	0
			2	2		
4	B	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
5	B	1	Total 27	C 10	N 5	O 10	P 2	0	0
5	C	1	Total 27	C 10	N 5	O 10	P 2	0	0
5	D	1	Total 27	C 10	N 5	O 10	P 2	0	0
5	E	1	Total 27	C 10	N 5	O 10	P 2	0	0

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 6 3 3	0	0
6	B	1	Total C O 6 3 3	0	0
6	C	1	Total C O 6 3 3	0	0
6	C	1	Total C O 6 3 3	0	0
6	D	1	Total C O 6 3 3	0	0
6	E	1	Total C O 6 3 3	0	0

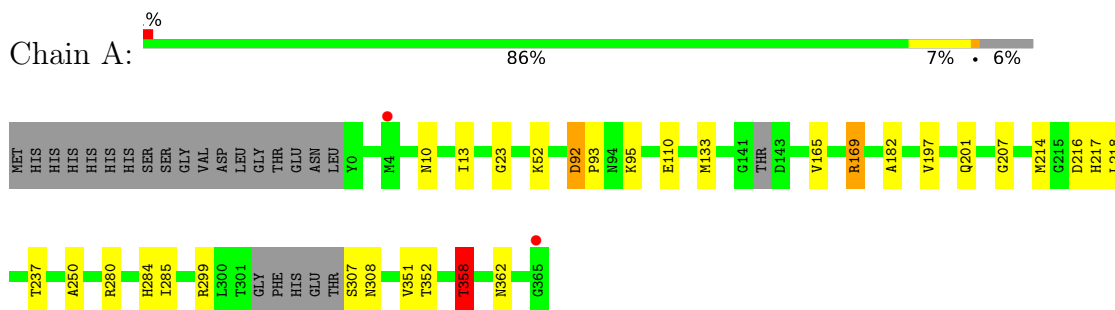
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	292	Total O 292 292	0	0
7	B	286	Total O 286 286	0	0
7	C	275	Total O 275 275	0	0
7	D	273	Total O 273 273	0	0
7	E	264	Total O 264 264	0	0

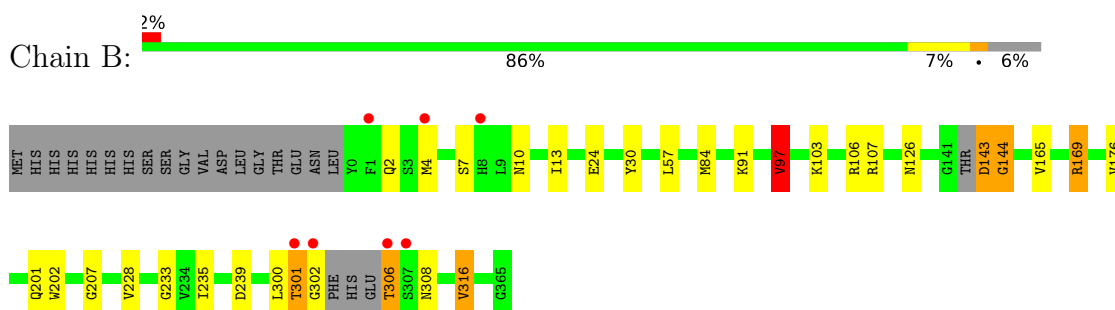
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

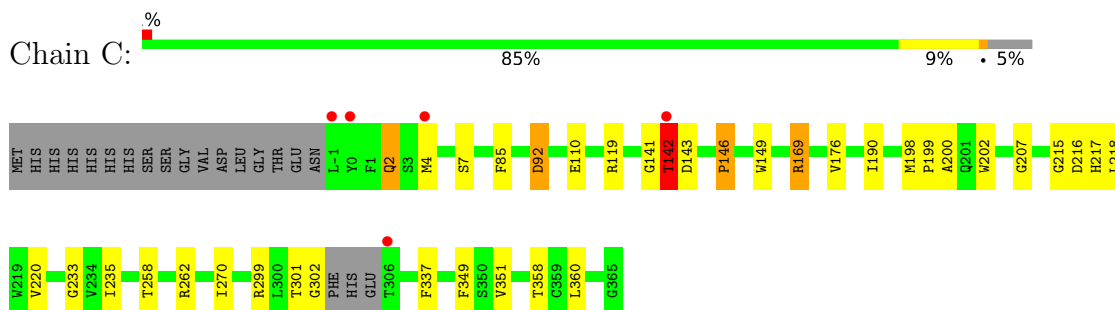
- Molecule 1: Glutamine synthetase



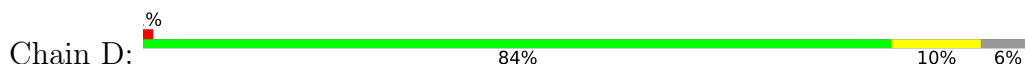
- Molecule 1: Glutamine synthetase

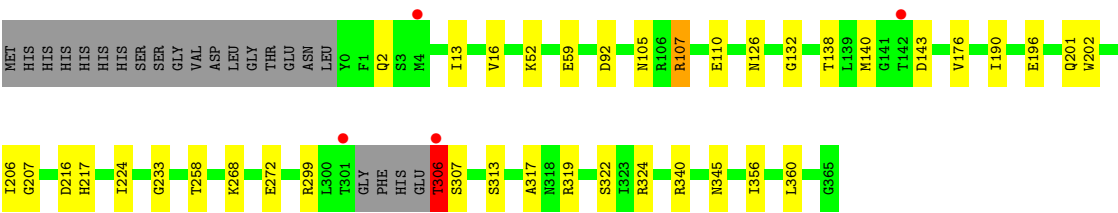


- Molecule 1: Glutamine synthetase

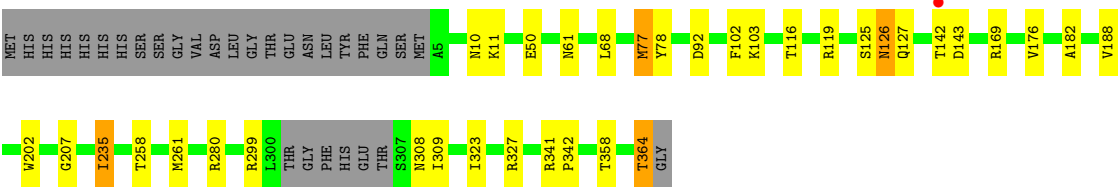
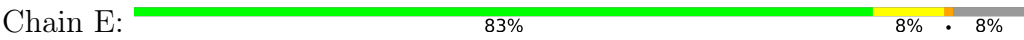


- Molecule 1: Glutamine synthetase





● Molecule 1: Glutamine synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.60Å 122.60Å 126.60Å 90.00° 130.60° 90.00°	Depositor
Resolution (Å)	39.25 – 2.05 39.25 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.25-2.05) 99.2 (39.25-2.05)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.3.0021	Depositor
R, R_{free}	0.159 , 0.212 0.165 , 0.213	Depositor DCC
R_{free} test set	6388 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15882	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, GOL, MN, CL, PO4

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	1.19	7/2932 (0.2%)	1.10	4/3961 (0.1%)
1	B	1.13	2/2947 (0.1%)	1.06	4/3981 (0.1%)
1	C	1.18	7/2949 (0.2%)	1.11	8/3987 (0.2%)
1	D	1.15	5/2938 (0.2%)	1.09	0/3972
1	E	1.13	3/2885 (0.1%)	1.08	4/3900 (0.1%)
All	All	1.16	24/14651 (0.2%)	1.09	20/19801 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	B	0	4
1	C	1	4
1	D	0	5
1	E	0	3
All	All	1	21

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	224	ILE	CA-CB	7.89	1.63	1.54
1	E	188	VAL	CA-CB	7.10	1.62	1.54
1	A	216	ASP	C-O	-6.81	1.16	1.24
1	D	217	HIS	C-O	-6.67	1.16	1.24
1	C	217	HIS	C-O	-6.63	1.16	1.24
1	A	218	LEU	C-O	-6.39	1.16	1.24
1	A	217	HIS	C-O	-6.33	1.16	1.24
1	A	237	THR	CA-CB	5.98	1.62	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	285	ILE	CA-CB	5.91	1.60	1.54
1	C	216	ASP	C-O	-5.59	1.17	1.24
1	B	316	VAL	CA-CB	5.57	1.60	1.54
1	E	142	THR	CA-C	5.55	1.60	1.52
1	C	351	VAL	CA-CB	5.50	1.60	1.54
1	B	97	VAL	CA-CB	5.48	1.61	1.54
1	A	351	VAL	CA-CB	5.39	1.60	1.54
1	D	190	ILE	CA-CB	5.35	1.60	1.54
1	D	206	ILE	CA-CB	5.35	1.60	1.54
1	C	169	ARG	CA-C	5.33	1.57	1.52
1	E	182	ALA	CA-CB	5.26	1.61	1.53
1	D	216	ASP	C-O	-5.24	1.18	1.24
1	C	146	PRO	CA-C	5.22	1.58	1.52
1	A	182	ALA	CA-CB	5.17	1.61	1.53
1	C	190	ILE	CA-CB	5.14	1.60	1.54
1	C	141	GLY	N-CA	-5.11	1.38	1.45

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	50	GLU	CA-C-N	-7.55	112.22	119.85
1	E	50	GLU	C-N-CA	-7.55	112.22	119.85
1	C	142	THR	OG1-CB-CG2	7.25	123.80	109.30
1	A	197	VAL	N-CA-C	6.79	117.58	110.72
1	A	308	ASN	N-CA-C	-6.70	100.52	110.23
1	B	144	GLY	N-CA-C	6.21	127.89	113.18
1	C	92	ASP	N-CA-C	6.15	117.13	109.57
1	C	262	ARG	NE-CZ-NH2	-6.14	113.67	119.20
1	B	228	VAL	CB-CA-C	-6.10	104.16	111.97
1	C	142	THR	N-CA-C	-5.80	105.30	112.90
1	A	358	THR	N-CA-CB	5.79	118.72	110.16
1	C	358	THR	N-CA-C	5.71	117.31	111.14
1	B	103	LYS	N-CA-C	5.46	117.12	110.41
1	C	92	ASP	CB-CA-C	-5.39	101.26	109.50
1	E	143	ASP	N-CA-C	5.32	122.13	110.80
1	A	92	ASP	N-CA-C	5.21	115.97	109.57
1	C	218	LEU	N-CA-C	5.20	116.64	111.07
1	C	142	THR	CA-CB-CG2	5.18	119.31	110.50
1	B	239	ASP	N-CA-C	5.13	115.88	109.57
1	E	323	ILE	CB-CA-C	-5.12	103.05	110.63

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	142	THR	CB

All (21) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	207	GLY	Mainchain,Peptide
1	A	307	SER	Peptide
1	A	92	ASP	Mainchain,Peptide
1	B	143	ASP	Peptide
1	B	207	GLY	Mainchain,Peptide
1	B	300	LEU	Peptide
1	C	207	GLY	Mainchain,Peptide
1	C	92	ASP	Mainchain,Peptide
1	D	207	GLY	Mainchain,Peptide
1	D	306	THR	Peptide
1	D	92	ASP	Mainchain,Peptide
1	E	207	GLY	Mainchain,Peptide
1	E	92	ASP	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2855	0	2743	17	1
1	B	2870	0	2760	24	0
1	C	2874	0	2757	26	0
1	D	2863	0	2747	27	0
1	E	2809	0	2705	20	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0
2	E	4	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
3	E	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	2	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	27	0	12	0	0
5	B	27	0	12	0	0
5	C	27	0	12	0	0
5	D	27	0	12	1	0
5	E	27	0	12	0	0
6	A	6	0	8	0	0
6	B	6	0	8	0	0
6	C	12	0	16	0	0
6	D	6	0	8	0	0
6	E	6	0	8	1	0
7	A	292	0	0	4	0
7	B	286	0	0	4	0
7	C	275	0	0	5	0
7	D	273	0	0	5	0
7	E	264	0	0	6	0
All	All	15882	0	13820	113	1

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 4.

All (113) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:LEU:HD11	1:B:84:MET:HE3	1.13	1.07
1:B:57:LEU:HD11	1:B:84:MET:CE	1.84	1.07
1:B:57:LEU:CD1	1:B:84:MET:HE3	1.84	1.07
1:C:2:GLN:HB3	1:C:235:ILE:HD11	1.42	0.99
1:D:138:THR:HG22	1:D:140:MET:HE2	1.46	0.97
1:C:4:MET:SD	1:C:142:THR:HG21	2.11	0.89
1:E:126:ASN:ND2	7:E:1660:HOH:O	2.10	0.85
1:C:301:THR:HG21	7:C:1702:HOH:O	1.77	0.83
1:C:110:GLU:OE1	7:C:1729:HOH:O	1.99	0.80
1:B:306:THR:N	7:B:1717:HOH:O	2.16	0.77
1:A:23:GLY:O	7:A:1659:HOH:O	2.04	0.75
1:D:138:THR:CG2	1:D:140:MET:HE2	2.15	0.75
1:D:59:GLU:OE2	7:D:1732:HOH:O	2.06	0.74
1:D:138:THR:CG2	1:D:140:MET:CE	2.67	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133[A]:MET:HE1	1:A:352:THR:HG23	1.71	0.71
1:D:306:THR:HG23	7:D:1679:HOH:O	1.93	0.69
1:B:126:ASN:OD1	7:B:1771:HOH:O	2.10	0.69
1:C:2:GLN:HB3	1:C:235:ILE:CD1	2.23	0.67
1:C:119:ARG:NE	7:C:1623:HOH:O	2.20	0.67
1:D:105:ASN:OD1	1:D:107:ARG:HB2	1.98	0.65
1:C:2:GLN:CG	1:C:142:THR:HA	2.28	0.63
1:E:127:GLN:OE1	7:E:1583:HOH:O	2.16	0.63
1:A:201:GLN:NE2	7:A:1594:HOH:O	2.31	0.63
1:B:2:GLN:HB3	1:B:235:ILE:HD11	1.81	0.62
1:D:317:ALA:HA	1:D:324:ARG:HD2	1.83	0.61
1:D:299:ARG:HD2	7:D:1682:HOH:O	2.00	0.61
1:A:284:HIS:NE2	1:A:358:THR:HG21	2.16	0.61
1:B:2:GLN:HB3	1:B:235:ILE:CD1	2.30	0.60
1:D:138:THR:HG22	1:D:140:MET:CE	2.24	0.60
1:B:302:GLY:HA2	1:B:306:THR:O	2.02	0.60
1:A:280:ARG:HE	1:A:358:THR:HG22	1.67	0.59
1:B:201:GLN:NE2	7:B:1635:HOH:O	2.36	0.59
1:A:299:ARG:HD2	7:A:1693:HOH:O	2.03	0.58
1:D:313:SER:O	1:D:322:SER:HB2	2.04	0.58
1:A:284:HIS:NE2	1:A:358:THR:CG2	2.66	0.58
1:C:2:GLN:CD	1:C:142:THR:HA	2.28	0.58
1:D:306:THR:CG2	7:D:1679:HOH:O	2.50	0.57
1:A:133[A]:MET:HE3	1:A:214:MET:HG2	1.84	0.57
1:B:57:LEU:CD1	1:B:84:MET:CE	2.62	0.57
1:C:4:MET:SD	1:C:142:THR:CG2	2.90	0.57
1:A:133[A]:MET:HE3	1:A:214:MET:CG	2.35	0.56
1:E:280:ARG:HH11	1:E:280:ARG:HG2	1.69	0.56
1:A:133[A]:MET:HE1	1:A:352:THR:CG2	2.36	0.55
1:C:2:GLN:NE2	1:C:142:THR:HG22	2.22	0.54
1:E:235:ILE:HD12	7:E:1732:HOH:O	2.07	0.54
1:B:30:TYR:HA	1:B:97:VAL:HG22	1.88	0.53
1:D:268:LYS:O	1:D:272:GLU:HG3	2.08	0.53
1:B:308:ASN:OD1	1:B:308:ASN:C	2.52	0.53
1:A:165:VAL:O	1:A:169:ARG:HG3	2.09	0.53
7:D:1510:HOH:O	1:E:327:ARG:HD2	2.09	0.52
1:B:97:VAL:O	1:B:97:VAL:HG23	2.11	0.51
1:B:165:VAL:O	1:B:169:ARG:HG3	2.11	0.51
1:D:319:ARG:HG2	1:D:340:ARG:CZ	2.40	0.51
1:A:133[A]:MET:CE	1:A:352:THR:CG2	2.89	0.50
1:C:299:ARG:HD2	7:C:1730:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:299:ARG:HD2	7:E:1563:HOH:O	2.13	0.49
1:C:143:ASP:OD1	1:C:143:ASP:N	2.42	0.49
1:E:77:MET:HE3	1:E:103:LYS:HA	1.93	0.49
1:D:319:ARG:HG2	1:D:340:ARG:NH1	2.27	0.49
1:E:280:ARG:HD3	1:E:358:THR:HG23	1.95	0.48
1:C:149:TRP:CZ2	1:C:198:MET:HE3	2.49	0.48
1:D:16:VAL:HG12	6:E:1505:GOL:H12	1.95	0.47
1:E:116:THR:HG22	1:E:119:ARG:HH21	1.79	0.47
1:D:138:THR:HG21	1:D:140:MET:CE	2.42	0.47
1:A:280:ARG:HE	1:A:358:THR:CG2	2.27	0.47
1:E:78:TYR:HB2	1:E:102:PHE:HB2	1.97	0.47
1:C:2:GLN:CD	1:C:142:THR:HG22	2.41	0.46
1:C:85:PHE:CE2	1:C:220:VAL:HG21	2.51	0.46
1:C:176:VAL:HG21	1:C:202:TRP:CE3	2.51	0.46
1:E:116:THR:HG22	1:E:119:ARG:NH2	2.30	0.46
1:A:250:ALA:HB2	7:A:1763:HOH:O	2.16	0.45
1:B:176:VAL:HG21	1:B:202:TRP:CD2	2.52	0.45
1:C:169:ARG:NE	7:C:1578:HOH:O	2.49	0.45
1:E:176:VAL:HG21	1:E:202:TRP:CD2	2.52	0.45
1:D:306:THR:CG2	1:D:307:SER:H	2.30	0.45
1:E:176:VAL:HG21	1:E:202:TRP:CE3	2.52	0.45
1:D:306:THR:HG23	1:D:307:SER:H	1.82	0.44
1:E:308:ASN:OD1	1:E:309:ILE:N	2.50	0.44
1:B:10:ASN:HB3	1:B:13:ILE:HD12	2.00	0.44
1:C:199:PRO:O	1:C:200:ALA:HB3	2.16	0.44
1:E:341:ARG:O	1:E:342:PRO:C	2.61	0.44
1:D:138:THR:HG21	1:D:140:MET:HE1	2.00	0.43
1:D:176:VAL:HG21	1:D:202:TRP:CE3	2.52	0.43
1:E:61:ASN:ND2	7:E:1591:HOH:O	2.28	0.43
1:B:302:GLY:C	1:B:306:THR:HG22	2.43	0.43
1:A:284:HIS:NE2	1:A:358:THR:HG23	2.33	0.43
1:C:215:GLY:HA3	1:C:349:PHE:CE1	2.54	0.43
1:D:176:VAL:HG21	1:D:202:TRP:CD2	2.54	0.43
1:C:2:GLN:HG2	1:C:142:THR:HA	2.01	0.43
1:C:258:THR:HG21	1:C:360:LEU:HD21	2.01	0.43
1:E:341:ARG:N	1:E:342:PRO:CD	2.83	0.42
1:C:176:VAL:HG21	1:C:202:TRP:CD2	2.54	0.42
1:B:97:VAL:O	1:B:97:VAL:CG2	2.67	0.42
1:A:93:PRO:O	1:A:95:LYS:HE2	2.20	0.42
1:C:146:PRO:HB2	1:C:149:TRP:CD1	2.54	0.42
1:B:176:VAL:HG21	1:B:202:TRP:CE3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:301:THR:HG23	1:C:302:GLY:O	2.19	0.42
1:B:308:ASN:OD1	7:B:1545:HOH:O	2.22	0.42
1:D:196:GLU:HB3	1:D:201:GLN:HE21	1.85	0.42
1:A:10:ASN:HB3	1:A:13:ILE:HD12	2.01	0.42
1:B:302:GLY:CA	1:B:306:THR:O	2.66	0.42
1:B:7:SER:HB3	1:B:233:GLY:HA3	2.02	0.41
1:B:301:THR:O	1:B:306:THR:HG22	2.20	0.41
1:B:57:LEU:CG	1:B:84:MET:CE	2.98	0.41
1:D:2:GLN:NE2	1:D:233:GLY:O	2.53	0.41
1:D:143:ASP:OD1	1:D:143:ASP:N	2.53	0.41
1:C:7:SER:HB3	1:C:233:GLY:HA3	2.03	0.41
1:D:13:ILE:HD11	1:E:11:LYS:HG3	2.03	0.41
1:D:132:GLY:HA3	5:D:501:ADP:H1'	2.03	0.41
1:D:356:ILE:HD12	1:D:360:LEU:HD12	2.03	0.41
1:C:270:ILE:HG23	1:C:337:PHE:CD2	2.57	0.40
1:E:258:THR:OG1	1:E:261:MET:HG3	2.21	0.40
1:E:364:THR:HG22	7:E:1739:HOH:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:ASN:ND2	1:A:362:ASN:ND2[2_554]	1.92	0.28

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	355/384 (92%)	346 (98%)	9 (2%)	0	100	100
1	B	357/384 (93%)	344 (96%)	12 (3%)	1 (0%)	36	30
1	C	360/384 (94%)	349 (97%)	11 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	358/384 (93%)	345 (96%)	13 (4%)	0	100	100
1	E	351/384 (91%)	339 (97%)	12 (3%)	0	100	100
All	All	1781/1920 (93%)	1723 (97%)	57 (3%)	1 (0%)	48	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	144	GLY

5.3.2 Protein sidechains [i](#)

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	302/323 (94%)	298 (99%)	4 (1%)	61	63
1	B	304/323 (94%)	293 (96%)	11 (4%)	31	26
1	C	304/323 (94%)	302 (99%)	2 (1%)	76	80
1	D	303/323 (94%)	296 (98%)	7 (2%)	44	42
1	E	298/323 (92%)	290 (97%)	8 (3%)	39	36
All	All	1511/1615 (94%)	1479 (98%)	32 (2%)	47	46

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LYS
1	A	110	GLU
1	A	169	ARG
1	A	358	THR
1	B	4	MET
1	B	24	GLU
1	B	91	LYS
1	B	97	VAL
1	B	106	ARG
1	B	107	ARG

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Mol	Chain	Res	Type
1	B	143	ASP
1	B	169	ARG
1	B	301	THR
1	B	306	THR
1	B	316	VAL
1	C	2	GLN
1	C	142	THR
1	D	52	LYS
1	D	107	ARG
1	D	110	GLU
1	D	126	ASN
1	D	258	THR
1	D	306	THR
1	D	345	ASN
1	E	10	ASN
1	E	68	LEU
1	E	77	MET
1	E	125	SER
1	E	126	ASN
1	E	169	ARG
1	E	235	ILE
1	E	364	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	201	GLN
1	A	248	ASN
1	A	255	ASN
1	B	201	GLN
1	B	248	ASN
1	B	282	GLN
1	B	310	ASN
1	C	8	HIS
1	D	126	ASN
1	D	201	GLN
1	D	248	ASN
1	D	265	ASN
1	D	282	GLN
1	D	345	ASN
1	E	10	ASN

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Mol	Chain	Res	Type
1	E	127	GLN
1	E	310	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 41 ligands modelled in this entry, 25 are monoatomic - leaving 16 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
6	GOL	C	1503	-	5,5,5	0.67	0	5,5,5	0.71	0
5	ADP	B	501	2	28,29,29	1.68	5 (17%)	43,45,45	1.90	11 (25%)
6	GOL	A	1501	-	5,5,5	0.55	0	5,5,5	0.89	0
3	PO4	A	502	2	4,4,4	1.40	0	6,6,6	1.15	0
5	ADP	D	501	2	28,29,29	1.65	5 (17%)	43,45,45	2.03	11 (25%)
6	GOL	E	1505	-	5,5,5	0.74	0	5,5,5	0.84	0
3	PO4	B	502	2	4,4,4	1.08	0	6,6,6	1.04	1 (16%)
3	PO4	C	502	2	4,4,4	0.68	0	6,6,6	2.27	2 (33%)
3	PO4	D	502	2	4,4,4	1.23	0	6,6,6	1.55	1 (16%)
6	GOL	C	1506	-	5,5,5	0.87	0	5,5,5	0.78	0
6	GOL	D	1504	-	5,5,5	0.52	0	5,5,5	1.03	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PO4	E	502	2	4,4,4	1.21	0	6,6,6	1.04	0
5	ADP	C	501	2	28,29,29	1.57	4 (14%)	43,45,45	2.17	12 (27%)
6	GOL	B	1502	-	5,5,5	0.72	0	5,5,5	0.88	0
5	ADP	A	501	2	28,29,29	1.39	3 (10%)	43,45,45	1.95	13 (30%)
5	ADP	E	501	2	28,29,29	1.54	4 (14%)	43,45,45	1.93	13 (30%)

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
6	GOL	C	1503	-	-	0/4/4/4	-
5	ADP	B	501	2	-	6/16/32/32	0/3/3/3
6	GOL	A	1501	-	-	2/4/4/4	-
5	ADP	D	501	2	-	5/16/32/32	0/3/3/3
6	GOL	E	1505	-	-	0/4/4/4	-
6	GOL	C	1506	-	-	2/4/4/4	-
6	GOL	D	1504	-	-	2/4/4/4	-
5	ADP	C	501	2	-	7/16/32/32	0/3/3/3
6	GOL	B	1502	-	-	0/4/4/4	-
5	ADP	A	501	2	-	4/16/32/32	0/3/3/3
5	ADP	E	501	2	-	5/16/32/32	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	501	ADP	C5-C4	5.08	1.48	1.39
5	B	501	ADP	PA-O3A	4.96	1.64	1.59
5	C	501	ADP	PA-O3A	4.58	1.64	1.59
5	E	501	ADP	C5-C4	4.35	1.46	1.39
5	D	501	ADP	PA-O3A	4.26	1.64	1.59
5	C	501	ADP	C5-C4	4.16	1.46	1.39
5	A	501	ADP	C5-C4	4.06	1.46	1.39
5	B	501	ADP	C5-C4	3.86	1.46	1.39
5	C	501	ADP	C5-C6	3.59	1.51	1.41
5	D	501	ADP	C5-C6	3.18	1.49	1.41
5	E	501	ADP	PA-O3A	3.07	1.62	1.59
5	E	501	ADP	C5-C6	3.04	1.49	1.41
5	C	501	ADP	C8-N7	2.79	1.37	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	501	ADP	C8-N7	2.76	1.37	1.31
5	E	501	ADP	C8-N7	2.74	1.36	1.31
5	D	501	ADP	C4-N9	-2.61	1.32	1.37
5	A	501	ADP	PB-O2B	-2.44	1.45	1.54
5	A	501	ADP	C5-C6	2.40	1.47	1.41
5	B	501	ADP	C5-N7	-2.25	1.35	1.39
5	B	501	ADP	C4-N9	-2.23	1.33	1.37
5	D	501	ADP	C8-N7	2.10	1.35	1.31

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	501	ADP	C5-C4-N3	-6.14	118.26	126.72
5	E	501	ADP	C5-C4-N3	-5.15	119.62	126.72
5	D	501	ADP	C5-C4-N3	-5.13	119.65	126.72
5	D	501	ADP	N3-C4-N9	5.07	135.79	127.17
5	B	501	ADP	C5-C4-N3	-5.02	119.80	126.72
5	D	501	ADP	N3-C2-N1	-4.97	121.07	128.58
5	B	501	ADP	N3-C4-N9	4.66	135.10	127.17
5	C	501	ADP	C2-N3-C4	4.65	123.18	111.83
5	A	501	ADP	N3-C2-N1	-4.63	121.58	128.58
5	C	501	ADP	N3-C4-N9	4.60	134.99	127.17
5	E	501	ADP	N3-C4-N9	4.54	134.89	127.17
5	C	501	ADP	N3-C2-N1	-4.49	121.78	128.58
5	D	501	ADP	C2-N3-C4	4.39	122.54	111.83
5	A	501	ADP	C5-C4-N3	-4.32	120.76	126.72
5	B	501	ADP	N3-C2-N1	-4.26	122.13	128.58
5	A	501	ADP	N3-C4-N9	4.16	134.25	127.17
3	C	502	PO4	O4-P-O1	-4.07	96.56	110.95
5	E	501	ADP	C4-N9-C8	3.99	109.93	105.74
5	B	501	ADP	C4-N9-C8	3.96	109.89	105.74
5	C	501	ADP	C4-C5-N7	-3.94	106.08	110.58
5	E	501	ADP	N3-C2-N1	-3.89	122.69	128.58
5	E	501	ADP	C2-N3-C4	3.70	120.87	111.83
5	D	501	ADP	C4-N9-C8	3.52	109.44	105.74
5	B	501	ADP	C2-N3-C4	3.47	120.31	111.83
5	A	501	ADP	C2-N3-C4	3.45	120.26	111.83
5	C	501	ADP	N9-C8-N7	-3.41	109.10	113.94
5	C	501	ADP	C5-N7-C8	3.40	108.80	103.45
5	C	501	ADP	C4-N9-C8	3.36	109.26	105.74
5	A	501	ADP	C4-N9-C8	3.25	109.16	105.74
5	E	501	ADP	N9-C8-N7	-3.15	109.47	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	501	ADP	O3'-C3'-C2'	-2.98	102.25	111.82
5	C	501	ADP	O4'-C1'-N9	2.95	113.75	108.09
5	C	501	ADP	C6-C5-N7	2.94	137.77	132.09
5	B	501	ADP	O3B-PB-O2B	2.88	118.59	107.80
5	E	501	ADP	C4-C5-N7	-2.81	107.37	110.58
5	A	501	ADP	C2-N1-C6	2.77	123.28	118.73
5	E	501	ADP	C5-N7-C8	2.74	107.75	103.45
3	D	502	PO4	O4-P-O1	-2.73	101.28	110.95
5	A	501	ADP	N6-C6-N1	2.65	124.29	118.38
3	C	502	PO4	O4-P-O2	2.63	116.09	107.91
5	D	501	ADP	C2-N1-C6	2.60	123.00	118.73
5	B	501	ADP	C2-N1-C6	2.55	122.92	118.73
5	E	501	ADP	C4-N9-C1'	-2.53	120.72	126.63
5	B	501	ADP	N6-C6-N1	2.52	123.99	118.38
5	A	501	ADP	C4-N9-C1'	-2.40	121.02	126.63
5	D	501	ADP	N6-C6-N1	2.39	123.69	118.38
5	B	501	ADP	N9-C8-N7	-2.34	110.62	113.94
5	D	501	ADP	C5-C6-N6	-2.31	117.56	123.29
5	A	501	ADP	C5-C6-N6	-2.31	117.56	123.29
5	E	501	ADP	O2B-PB-O1B	2.25	119.58	110.83
5	E	501	ADP	C2-N1-C6	2.21	122.36	118.73
3	B	502	PO4	O3-P-O2	2.19	114.72	107.91
5	C	501	ADP	O2B-PB-O1B	2.13	119.14	110.83
5	B	501	ADP	C5-C6-N6	-2.13	118.02	123.29
5	A	501	ADP	O2B-PB-O1B	2.12	119.11	110.83
5	A	501	ADP	O3'-C3'-C2'	-2.12	105.02	111.82
5	A	501	ADP	C4-C5-N7	-2.11	108.17	110.58
5	D	501	ADP	O2A-PA-O1A	2.10	122.21	112.44
5	B	501	ADP	O4'-C1'-N9	2.07	112.06	108.09
5	C	501	ADP	C4-N9-C1'	-2.05	121.83	126.63
5	E	501	ADP	N6-C6-N1	2.05	122.95	118.38
5	A	501	ADP	O2B-PB-O3A	-2.03	97.81	104.64
5	D	501	ADP	C4-N9-C1'	-2.02	121.91	126.63
5	E	501	ADP	C6-C5-N7	2.02	135.97	132.09

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	501	ADP	PA-O3A-PB-O3B
5	C	501	ADP	PA-O3A-PB-O3B
5	D	501	ADP	PA-O3A-PB-O3B

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Mol	Chain	Res	Type	Atoms
5	E	501	ADP	PA-O3A-PB-O3B
5	E	501	ADP	C5'-O5'-PA-O1A
5	D	501	ADP	PA-O3A-PB-O1B
5	C	501	ADP	O4'-C4'-C5'-O5'
5	A	501	ADP	PA-O3A-PB-O2B
5	E	501	ADP	C2'-C1'-N9-C8
6	A	1501	GOL	O1-C1-C2-O2
6	D	1504	GOL	O1-C1-C2-O2
5	A	501	ADP	C5'-O5'-PA-O1A
5	C	501	ADP	C2'-C1'-N9-C8
5	A	501	ADP	C2'-C1'-N9-C8
5	D	501	ADP	C2'-C1'-N9-C8
6	A	1501	GOL	O1-C1-C2-C3
6	D	1504	GOL	O1-C1-C2-C3
5	C	501	ADP	C3'-C4'-C5'-O5'
6	C	1506	GOL	O2-C2-C3-O3
5	B	501	ADP	O4'-C4'-C5'-O5'
5	D	501	ADP	PB-O3A-PA-O2A
5	C	501	ADP	O4'-C1'-N9-C8
5	A	501	ADP	PA-O3A-PB-O1B
5	B	501	ADP	PA-O3A-PB-O1B
5	C	501	ADP	PA-O3A-PB-O1B
5	E	501	ADP	PA-O3A-PB-O1B
5	B	501	ADP	PA-O3A-PB-O2B
5	B	501	ADP	C2'-C1'-N9-C8
5	C	501	ADP	PB-O3A-PA-O1A
6	C	1506	GOL	C1-C2-C3-O3
5	B	501	ADP	O4'-C1'-N9-C8
5	E	501	ADP	O4'-C1'-N9-C8
5	D	501	ADP	PB-O3A-PA-O1A

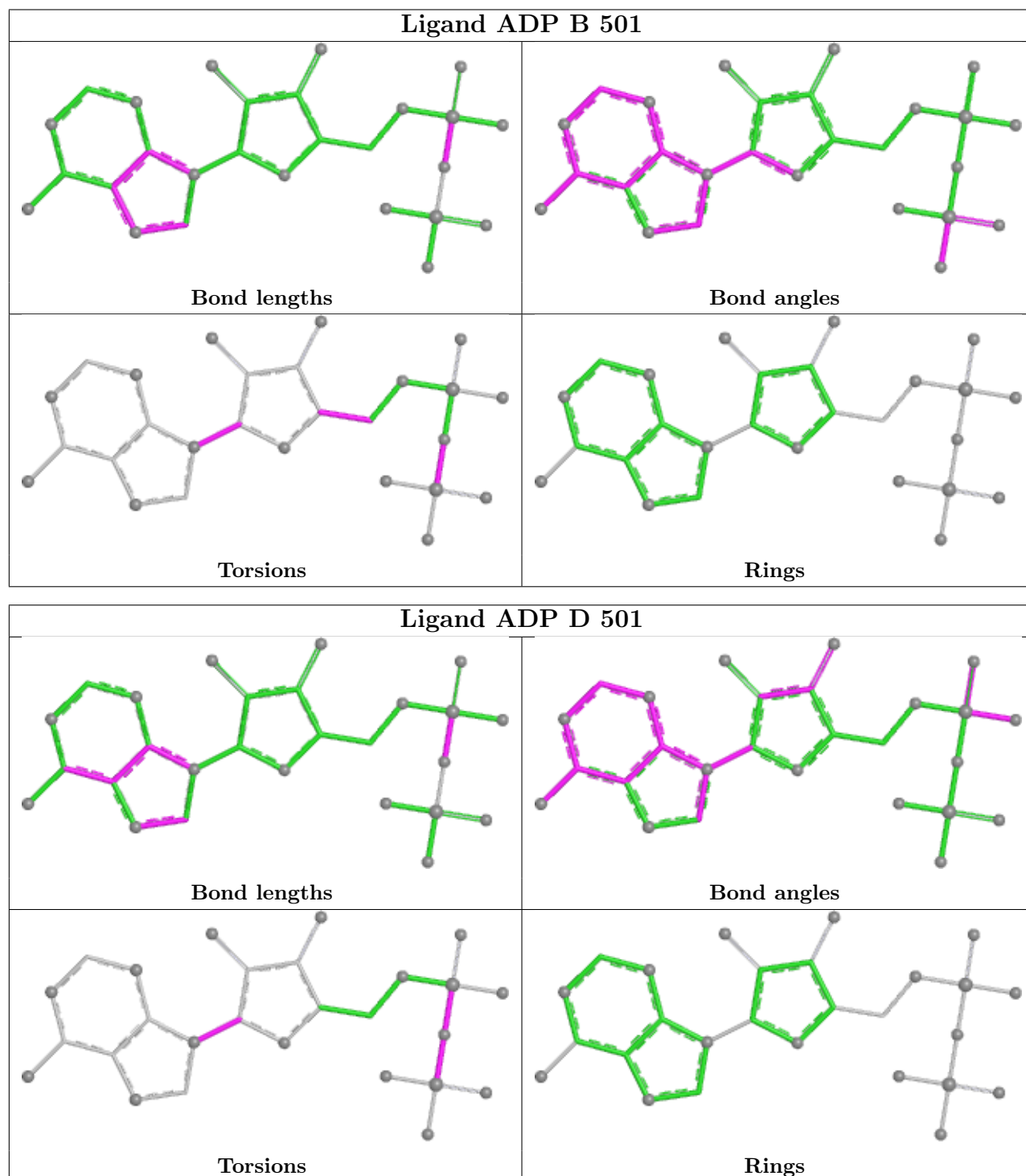
There are no ring outliers.

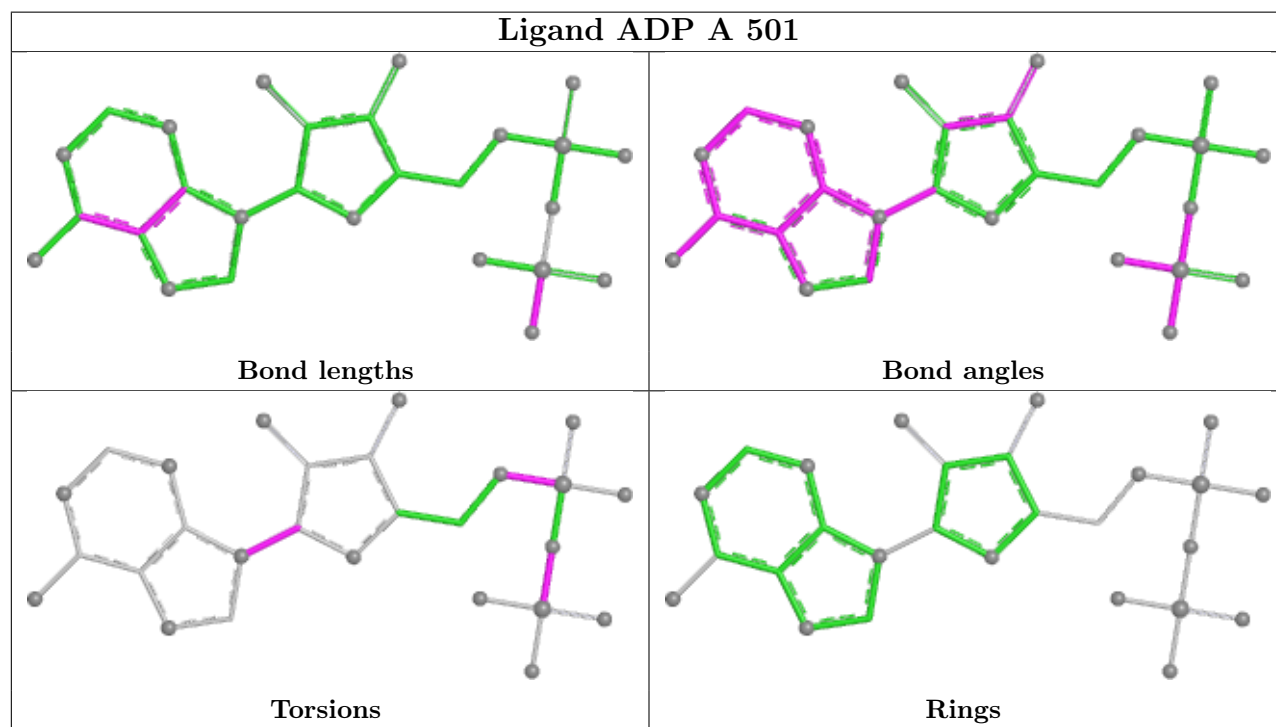
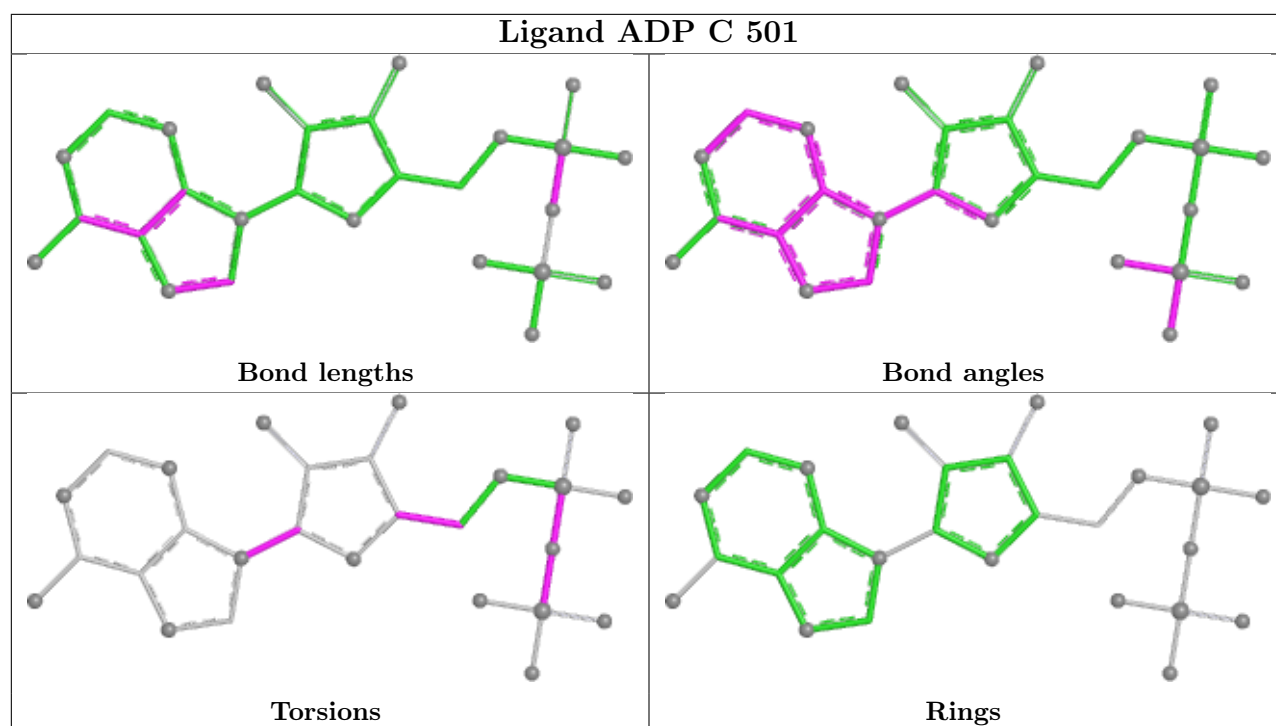
2 monomers are involved in 2 short contacts:

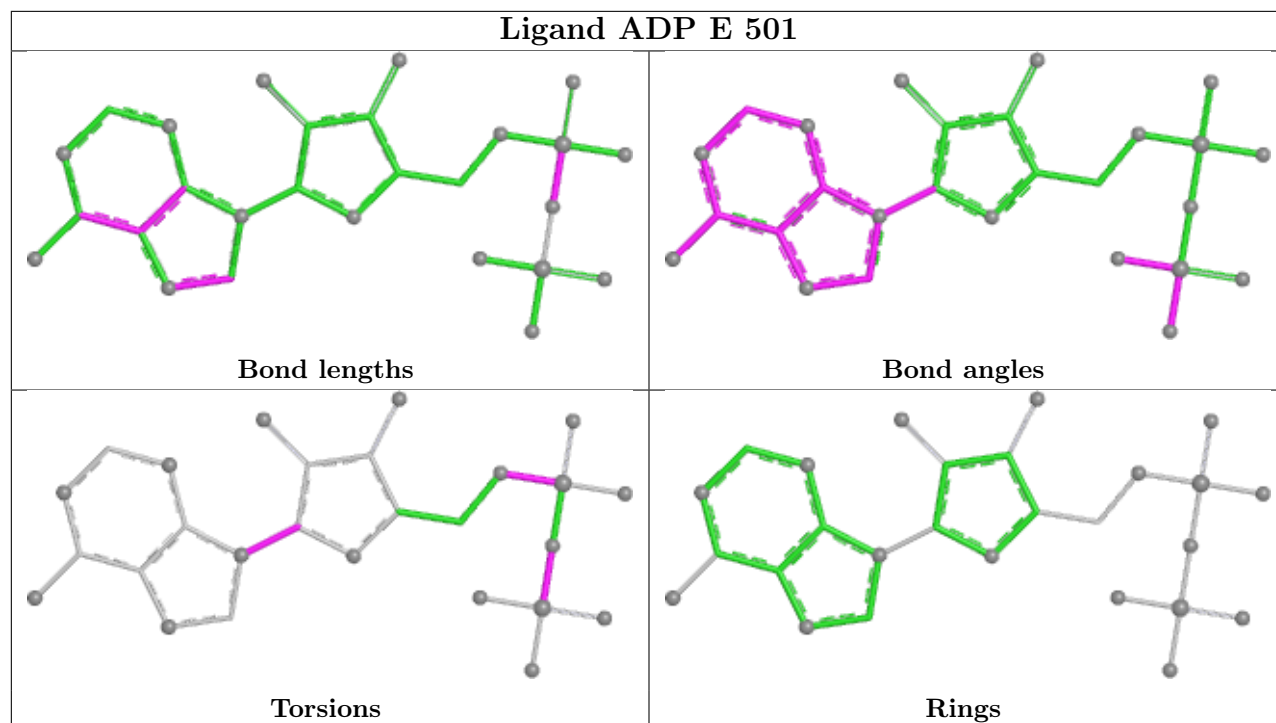
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	501	ADP	1	0
6	E	1505	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	360/384 (93%)	-0.54	2 (0%) 85 87	4, 12, 26, 33	1 (0%)
1	B	362/384 (94%)	-0.41	7 (1%) 66 68	5, 12, 27, 37	1 (0%)
1	C	364/384 (94%)	-0.49	5 (1%) 73 75	5, 12, 27, 33	0
1	D	362/384 (94%)	-0.46	4 (1%) 78 80	5, 13, 25, 32	0
1	E	354/384 (92%)	-0.37	1 (0%) 90 92	6, 13, 24, 31	1 (0%)
All	All	1802/1920 (93%)	-0.46	19 (1%) 78 80	4, 12, 26, 37	3 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	306	THR	4.1
1	B	306	THR	3.7
1	C	142	THR	3.5
1	D	301	THR	2.9
1	A	4	MET	2.8
1	B	301	THR	2.8
1	A	365	GLY	2.7
1	D	142	THR	2.7
1	B	4	MET	2.6
1	E	142	THR	2.4
1	C	4	MET	2.3
1	D	306	THR	2.3
1	B	302	GLY	2.3
1	B	8	HIS	2.3
1	C	0	TYR	2.2
1	D	4	MET	2.2
1	C	-1	LEU	2.1
1	B	307	SER	2.1
1	B	1	PHE	2.1

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

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

6.3 Carbohydrates [i](#)

There are no 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.

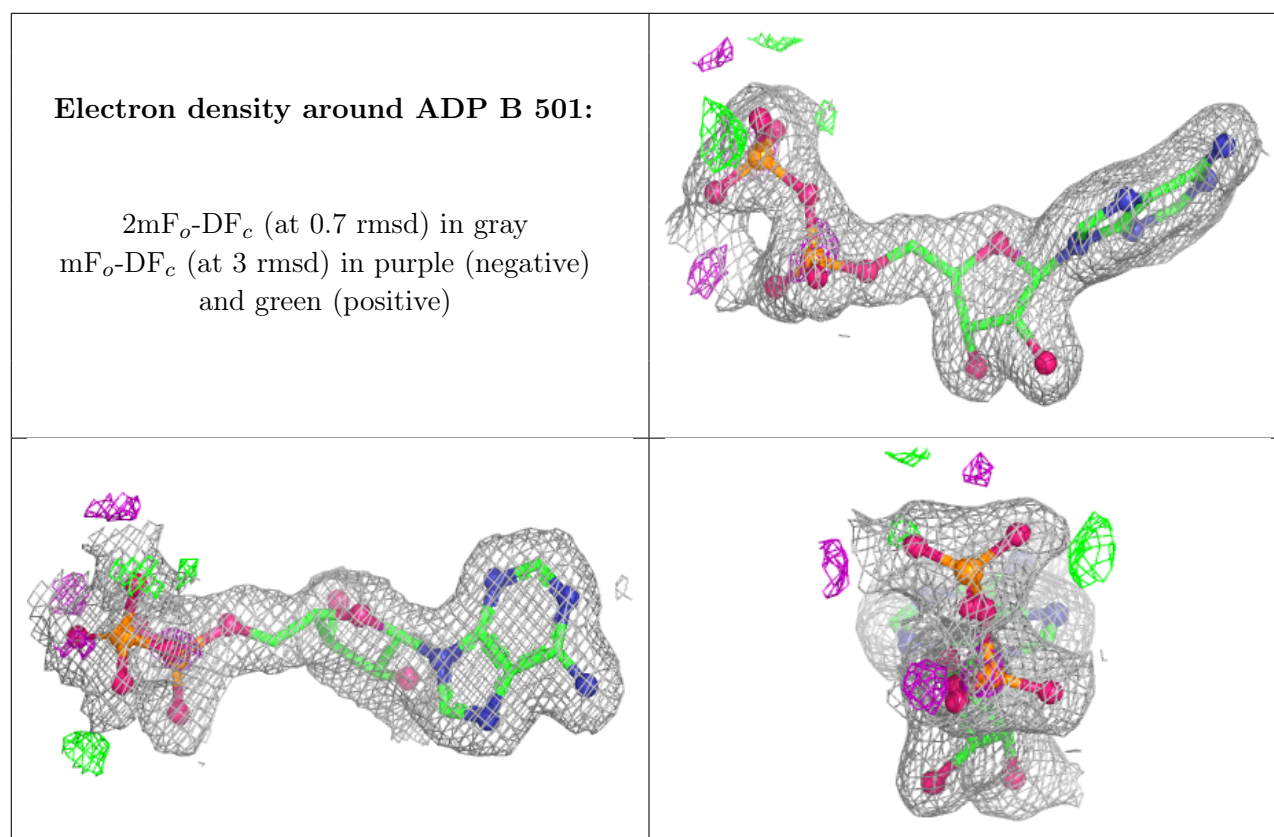
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	GOL	C	1506	6/6	0.90	0.09	24,28,29,29	0
6	GOL	C	1503	6/6	0.91	0.09	14,15,16,17	0
6	GOL	B	1502	6/6	0.91	0.10	14,16,19,19	0
6	GOL	D	1504	6/6	0.93	0.07	20,20,21,21	0
6	GOL	A	1501	6/6	0.94	0.11	16,18,18,19	0
6	GOL	E	1505	6/6	0.95	0.09	17,20,21,21	0
5	ADP	B	501	27/27	0.98	0.05	5,12,14,15	0
5	ADP	C	501	27/27	0.98	0.05	6,12,14,15	0
5	ADP	D	501	27/27	0.98	0.05	2,11,14,14	0
5	ADP	E	501	27/27	0.98	0.05	9,15,17,23	0
3	PO4	D	502	5/5	0.98	0.06	5,10,13,13	0
3	PO4	B	502	5/5	0.99	0.07	4,9,11,14	0
3	PO4	C	502	5/5	0.99	0.06	4,7,9,11	0
2	MN	A	402	1/1	0.99	0.07	8,8,8,8	0
3	PO4	E	502	5/5	0.99	0.06	9,10,12,13	0
4	CL	A	1403	1/1	0.99	0.05	9,9,9,9	0
4	CL	A	1405	1/1	0.99	0.07	9,9,9,9	0
4	CL	B	1401	1/1	0.99	0.05	10,10,10,10	0
4	CL	C	1402	1/1	0.99	0.07	9,9,9,9	0
4	CL	D	1404	1/1	0.99	0.08	10,10,10,10	0
5	ADP	A	501	27/27	0.99	0.04	4,9,14,15	0
2	MN	A	404	1/1	0.99	0.02	13,13,13,13	1
2	MN	B	403	1/1	0.99	0.04	8,8,8,8	0
2	MN	B	404	1/1	0.99	0.04	16,16,16,16	1
2	MN	C	402	1/1	0.99	0.05	10,10,10,10	0
2	MN	C	404	1/1	0.99	0.02	10,10,10,10	1
2	MN	D	402	1/1	0.99	0.05	11,11,11,11	0

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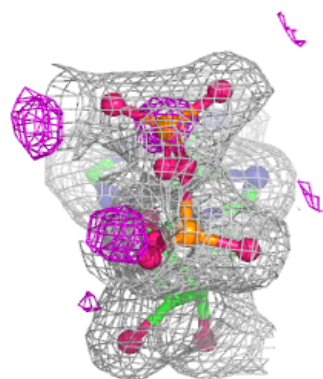
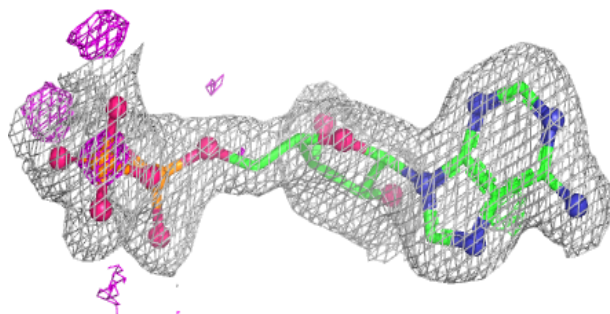
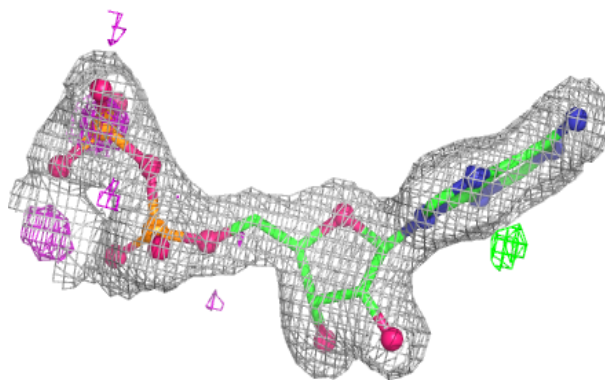
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MN	D	404	1/1	0.99	0.04	11,11,11,11	1
2	MN	E	402	1/1	0.99	0.03	13,13,13,13	0
2	MN	E	404	1/1	0.99	0.02	17,17,17,17	1
3	PO4	A	502	5/5	0.99	0.07	3,6,7,11	0
2	MN	B	402	1/1	1.00	0.05	10,10,10,10	0
2	MN	D	401	1/1	1.00	0.04	8,8,8,8	0
2	MN	A	403	1/1	1.00	0.05	6,6,6,6	0
2	MN	D	403	1/1	1.00	0.04	7,7,7,7	0
2	MN	A	401	1/1	1.00	0.04	7,7,7,7	0
2	MN	E	401	1/1	1.00	0.03	11,11,11,11	0
2	MN	C	401	1/1	1.00	0.05	11,11,11,11	0
2	MN	E	403	1/1	1.00	0.04	9,9,9,9	0
2	MN	B	401	1/1	1.00	0.05	9,9,9,9	0
2	MN	C	403	1/1	1.00	0.05	6,6,6,6	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.

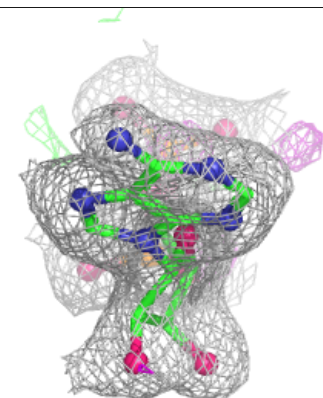
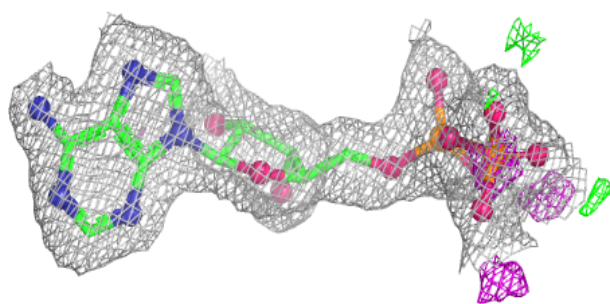
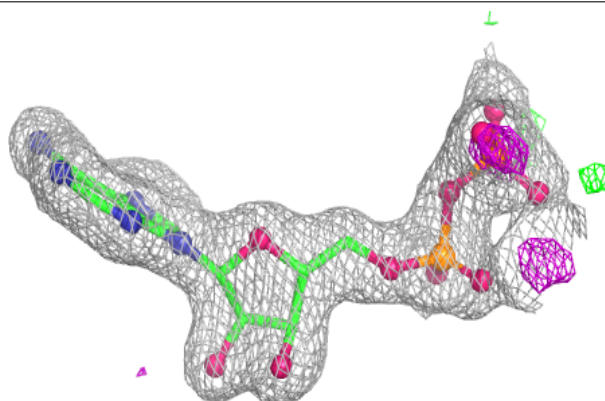


Electron density around ADP C 501:

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

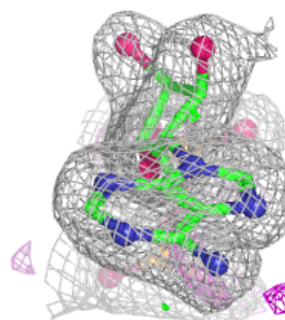
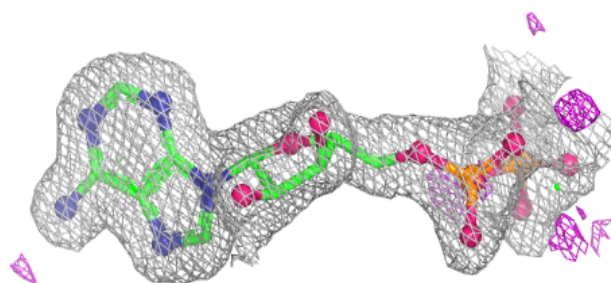
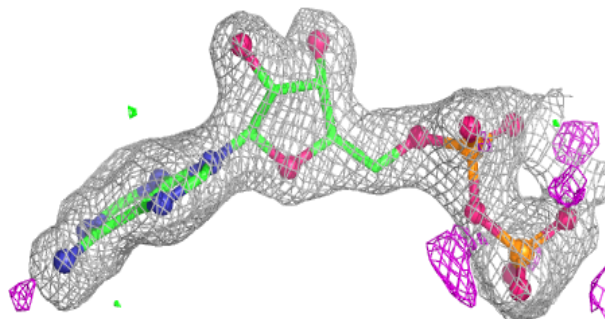
**Electron density around ADP D 501:**

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

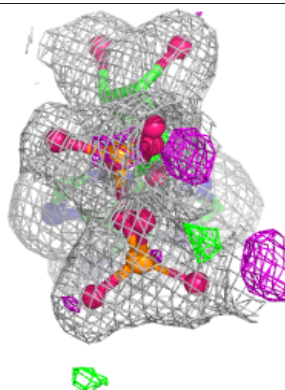
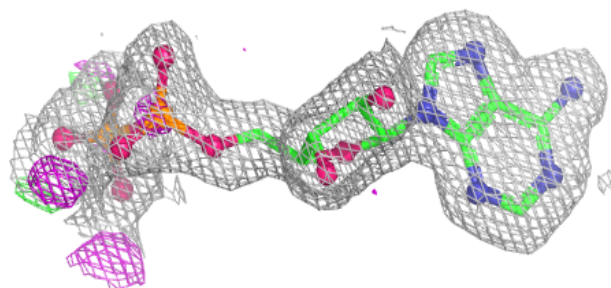
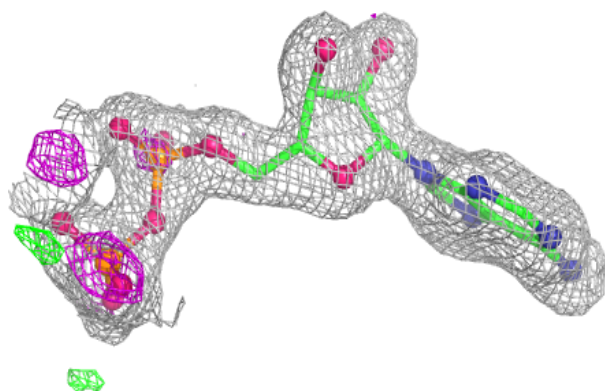


Electron density around ADP E 501:

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

**Electron density around ADP A 501:**

$2mF_o-DF_c$ (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.