



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2026 – 11:51 PM UTC

PDB ID : 1XJQ / pdb_00001xjq
Title : ADP Complex OF HUMAN PAPS SYNTHETASE 1
Authors : Harjes, S.; Bayer, P.; Scheidig, A.J.
Deposited on : 2004-09-24
Resolution : 2.06 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

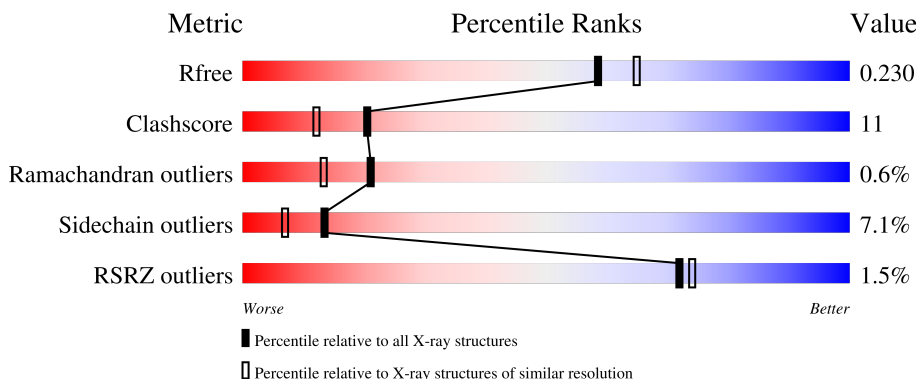
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	630	% 64% 20% 5% 11%
1	B	630	% 58% 28% 7% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9997 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 Bifunctional 3'-phosphoadenosine 5'-phosphosulfate synthetase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	590	Total	C	N	O	S	0	5	0
			4718	2990	830	867	31			
1	A	563	Total	C	N	O	S	0	8	0
			4514	2864	790	828	32			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	416	SER	PHE	conflict	UNP O43252
B	625	HIS	-	expression tag	UNP O43252
B	626	HIS	-	expression tag	UNP O43252
B	627	HIS	-	expression tag	UNP O43252
B	628	HIS	-	expression tag	UNP O43252
B	629	HIS	-	expression tag	UNP O43252
B	630	HIS	-	expression tag	UNP O43252
A	416	SER	PHE	conflict	UNP O43252
A	625	HIS	-	expression tag	UNP O43252
A	626	HIS	-	expression tag	UNP O43252
A	627	HIS	-	expression tag	UNP O43252
A	628	HIS	-	expression tag	UNP O43252
A	629	HIS	-	expression tag	UNP O43252
A	630	HIS	-	expression tag	UNP O43252

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



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

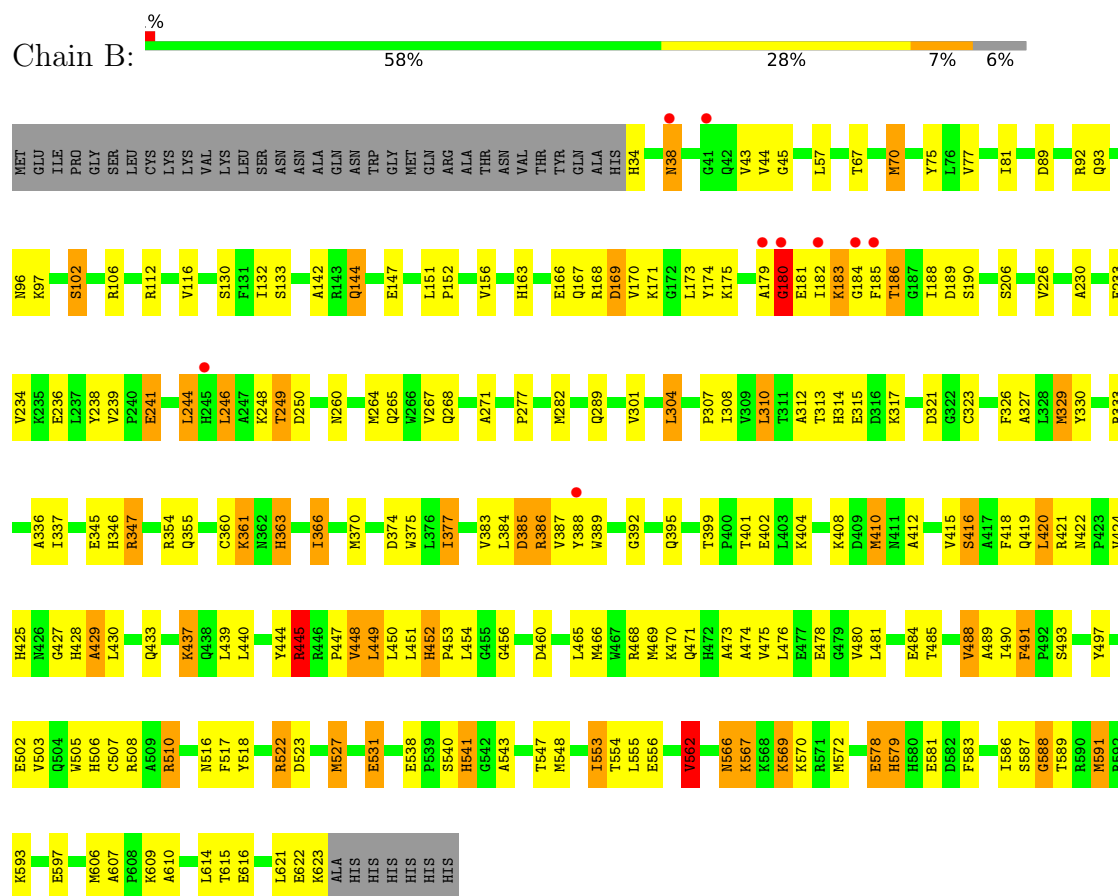
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	323	Total O 323 323	0	0
3	A	361	Total O 361 361	0	0

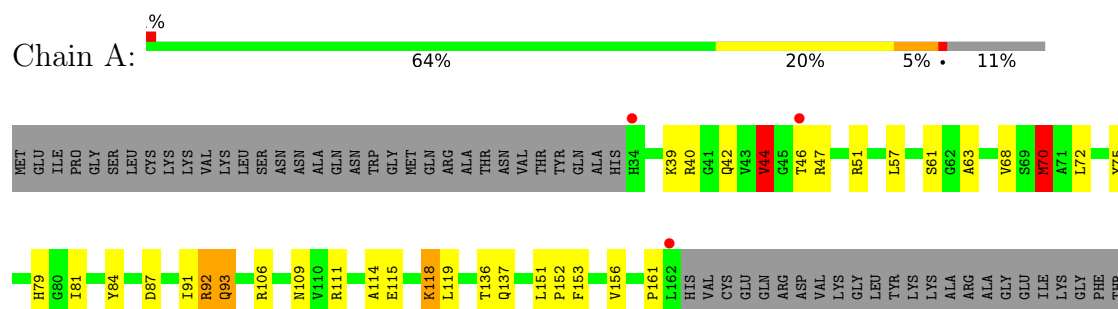
3 Residue-property plots

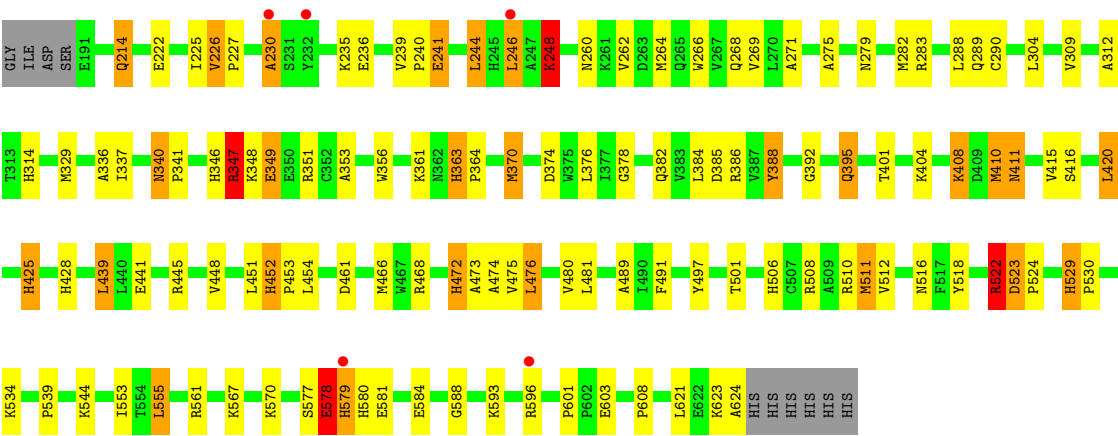
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Bifunctional 3'-phosphoadenosine 5'-phosphosulfate synthetase 1



- Molecule 1: Bifunctional 3'-phosphoadenosine 5'-phosphosulfate synthetase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.50Å 82.50Å 133.00Å 90.00° 105.00° 90.00°	Depositor
Resolution (Å)	129.10 – 2.06 128.47 – 2.06	Depositor EDS
% Data completeness (in resolution range)	96.6 (129.10-2.06) 96.5 (128.47-2.06)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.171 , 0.221 0.181 , 0.230	Depositor DCC
R_{free} test set	4869 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.304	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9997	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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

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	1.91	69/4657 (1.5%)	1.49	34/6314 (0.5%)
1	B	1.94	97/4854 (2.0%)	1.51	38/6574 (0.6%)
All	All	1.93	166/9511 (1.7%)	1.50	72/12888 (0.6%)

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	2
1	B	0	2
All	All	0	4

All (166) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	340	ASN	C-O	-16.05	1.16	1.23
1	A	370	MET	SD-CE	15.61	2.18	1.79
1	A	346	HIS	C-O	13.27	1.39	1.24
1	B	527	MET	SD-CE	12.07	2.09	1.79
1	B	70	MET	SD-CE	11.55	2.08	1.79
1	B	451	LEU	CA-C	-10.59	1.39	1.52
1	A	346	HIS	CA-C	10.38	1.65	1.52
1	A	70	MET	SD-CE	10.25	2.05	1.79
1	B	548	MET	SD-CE	-10.22	1.53	1.79
1	A	410	MET	SD-CE	-10.02	1.54	1.79
1	A	262	VAL	CA-CB	9.63	1.66	1.54
1	B	490	ILE	CA-C	9.20	1.63	1.52
1	B	591	MET	SD-CE	9.19	2.02	1.79
1	A	466	MET	SD-CE	8.64	2.01	1.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	448	VAL	C-O	8.63	1.32	1.24
1	B	236	GLU	CA-C	-8.51	1.41	1.52
1	B	420	LEU	CA-C	-8.41	1.42	1.52
1	B	491	PHE	CA-C	-8.34	1.43	1.52
1	B	81	ILE	CA-CB	-8.05	1.47	1.54
1	B	449	LEU	CA-C	-7.84	1.43	1.52
1	B	308	ILE	CA-C	-7.79	1.46	1.53
1	B	517	PHE	C-O	-7.67	1.14	1.23
1	A	244	LEU	C-O	7.51	1.32	1.24
1	B	470	LYS	N-CA	7.50	1.55	1.46
1	B	523	ASP	CA-CB	7.37	1.59	1.52
1	A	340	ASN	CA-C	-7.33	1.48	1.53
1	B	330	TYR	CA-C	7.27	1.61	1.52
1	A	336	ALA	CA-C	-7.27	1.43	1.52
1	B	265	GLN	CA-C	7.25	1.62	1.52
1	B	488	VAL	N-CA	-7.25	1.37	1.46
1	A	260	ASN	CA-C	-7.22	1.42	1.53
1	A	415	VAL	CA-CB	-7.21	1.45	1.54
1	A	410	MET	CG-SD	7.20	1.98	1.80
1	A	40	ARG	CG-CD	7.11	1.73	1.52
1	B	503	VAL	CA-CB	-7.08	1.44	1.54
1	B	505	TRP	CA-C	7.06	1.61	1.52
1	B	489	ALA	C-O	-7.05	1.15	1.23
1	B	469	MET	CA-C	7.01	1.61	1.52
1	A	269	VAL	CA-CB	6.97	1.62	1.54
1	A	44	VAL	C-O	6.90	1.31	1.24
1	B	469	MET	C-O	6.90	1.32	1.24
1	B	402	GLU	C-O	-6.89	1.16	1.24
1	A	518	TYR	C-O	-6.88	1.16	1.24
1	A	282	MET	SD-CE	-6.80	1.62	1.79
1	A	512	VAL	CA-CB	-6.78	1.45	1.54
1	A	466	MET	N-CA	6.77	1.54	1.46
1	A	111	ARG	NE-CZ	-6.73	1.25	1.33
1	A	226	VAL	CA-C	6.69	1.58	1.52
1	B	377	ILE	C-O	-6.62	1.17	1.24
1	A	269	VAL	C-O	-6.61	1.16	1.24
1	B	385	ASP	CA-C	-6.53	1.44	1.52
1	A	92	ARG	NE-CZ	6.52	1.40	1.33
1	B	329	MET	SD-CE	6.51	1.95	1.79
1	B	347	ARG	CZ-NH1	6.51	1.41	1.32
1	B	206	SER	C-O	-6.49	1.14	1.24
1	A	63	ALA	C-O	-6.48	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	337	ILE	CA-C	6.46	1.60	1.52
1	B	386	ARG	N-CA	-6.45	1.38	1.46
1	B	226	VAL	CA-CB	6.43	1.62	1.54
1	A	309	VAL	CA-CB	6.41	1.66	1.55
1	B	466	MET	SD-CE	6.37	1.95	1.79
1	B	327	ALA	CA-CB	6.37	1.63	1.53
1	B	449	LEU	C-O	6.35	1.31	1.24
1	B	437	LYS	CA-C	-6.31	1.44	1.52
1	A	401	THR	CB-CG2	6.28	1.73	1.52
1	B	454	LEU	CG-CD1	6.26	1.73	1.52
1	A	491	PHE	CA-C	-6.25	1.46	1.52
1	B	282	MET	SD-CE	-6.20	1.64	1.79
1	A	348	LYS	CA-C	-6.16	1.44	1.52
1	A	236	GLU	CA-C	-6.15	1.44	1.52
1	A	275	ALA	C-O	-6.15	1.15	1.24
1	B	418	PHE	CA-C	-6.14	1.45	1.53
1	B	67	THR	CA-C	6.10	1.60	1.52
1	B	412	ALA	CA-CB	6.09	1.63	1.53
1	A	448	VAL	CA-CB	-6.09	1.46	1.54
1	B	38	ASN	CB-CG	6.07	1.67	1.52
1	A	384	LEU	N-CA	6.06	1.54	1.46
1	A	475	VAL	CA-CB	5.97	1.61	1.54
1	B	453	PRO	CA-CB	5.96	1.61	1.53
1	B	355	GLN	CA-C	-5.90	1.45	1.52
1	B	327	ALA	CA-C	-5.87	1.45	1.52
1	A	114	ALA	CA-CB	-5.85	1.44	1.53
1	B	326	PHE	CA-C	-5.85	1.45	1.52
1	B	366	ILE	CA-CB	5.84	1.61	1.54
1	A	351	ARG	CA-C	-5.84	1.45	1.52
1	B	346	HIS	C-O	5.80	1.31	1.23
1	B	416[A]	SER	CA-CB	-5.79	1.44	1.53
1	B	416[B]	SER	CA-CB	-5.79	1.44	1.53
1	A	476[A]	LEU	CA-C	5.78	1.60	1.52
1	B	336	ALA	C-O	5.76	1.30	1.23
1	B	453	PRO	CA-C	-5.75	1.45	1.52
1	A	561	ARG	NE-CZ	5.75	1.39	1.33
1	A	271	ALA	CA-C	-5.75	1.45	1.52
1	B	389	TRP	CA-C	5.73	1.59	1.52
1	B	133	SER	CA-CB	5.70	1.56	1.52
1	A	523	ASP	CA-CB	5.68	1.57	1.52
1	B	473	ALA	CA-CB	5.63	1.62	1.53
1	B	361	LYS	CG-CD	5.62	1.69	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	84	TYR	CA-C	-5.62	1.45	1.52
1	B	415	VAL	N-CA	-5.61	1.39	1.46
1	A	386	ARG	N-CA	-5.60	1.39	1.46
1	B	260	ASN	CA-C	-5.60	1.45	1.53
1	A	451	LEU	CA-C	-5.59	1.46	1.52
1	A	347	ARG	CA-C	-5.59	1.45	1.52
1	A	81	ILE	CA-CB	-5.59	1.48	1.53
1	B	614	LEU	C-O	5.54	1.30	1.24
1	A	349	GLU	C-O	-5.54	1.17	1.24
1	B	336	ALA	CA-C	-5.52	1.45	1.52
1	B	421	ARG	CZ-NH2	5.52	1.40	1.33
1	A	68	VAL	CA-CB	-5.52	1.48	1.54
1	B	307	PRO	CA-C	5.51	1.59	1.52
1	A	474	ALA	CA-C	-5.50	1.45	1.52
1	B	132	ILE	CA-CB	5.49	1.60	1.54
1	A	404	LYS	CA-C	5.49	1.60	1.52
1	B	466	MET	N-CA	5.48	1.52	1.46
1	A	70	MET	CG-SD	-5.48	1.67	1.80
1	A	439	LEU	CA-C	5.47	1.59	1.52
1	B	244	LEU	C-O	5.46	1.30	1.24
1	B	384	LEU	N-CA	5.44	1.53	1.46
1	A	337	ILE	CA-C	5.44	1.59	1.52
1	A	522	ARG	C-O	-5.42	1.17	1.24
1	B	553	ILE	CA-CB	5.41	1.61	1.54
1	B	410	MET	SD-CE	5.40	1.93	1.79
1	A	529	HIS	CA-C	-5.38	1.47	1.53
1	B	43	VAL	CA-CB	5.38	1.61	1.54
1	B	142	ALA	C-O	-5.35	1.17	1.24
1	B	271	ALA	CA-CB	5.34	1.61	1.53
1	A	260	ASN	C-N	-5.34	1.26	1.33
1	B	147	GLU	CA-C	5.30	1.59	1.52
1	A	353	ALA	CA-CB	5.30	1.61	1.53
1	B	130	SER	C-O	-5.29	1.16	1.23
1	B	230	ALA	CA-C	5.29	1.59	1.52
1	A	248	LYS	N-CA	5.29	1.53	1.46
1	A	473	ALA	CA-CB	5.28	1.61	1.53
1	B	465	LEU	CA-C	5.28	1.59	1.52
1	B	289	GLN	CA-C	-5.27	1.46	1.52
1	B	444	TYR	CA-C	5.26	1.59	1.52
1	B	267	VAL	CA-C	-5.25	1.45	1.52
1	B	301	VAL	C-O	-5.24	1.18	1.24
1	A	312	ALA	C-O	5.23	1.30	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	376	LEU	CA-C	-5.23	1.46	1.52
1	A	425	HIS	N-CA	-5.21	1.39	1.45
1	A	480	VAL	CA-C	5.19	1.59	1.52
1	B	480	VAL	CA-C	5.19	1.60	1.52
1	B	147	GLU	CD-OE2	5.16	1.35	1.25
1	A	522	ARG	NE-CZ	5.16	1.38	1.33
1	B	249	THR	CA-CB	-5.15	1.45	1.53
1	B	493	SER	N-CA	5.15	1.53	1.45
1	A	153	PHE	C-O	5.15	1.29	1.24
1	B	541	HIS	CA-CB	5.14	1.61	1.53
1	B	401	THR	CA-C	5.14	1.59	1.52
1	B	562	VAL	C-O	-5.11	1.18	1.24
1	B	346	HIS	CA-CB	-5.10	1.46	1.52
1	B	566	ASN	CA-C	5.10	1.59	1.52
1	B	497	TYR	CA-C	5.09	1.59	1.53
1	B	282	MET	CA-C	-5.08	1.46	1.52
1	A	534	LYS	CD-CE	5.08	1.67	1.52
1	B	97	LYS	N-CA	-5.08	1.40	1.46
1	A	225	ILE	CA-CB	5.06	1.60	1.55
1	B	304	LEU	CA-C	-5.06	1.48	1.53
1	A	461	ASP	CA-CB	5.06	1.61	1.53
1	B	170	VAL	CA-CB	5.03	1.62	1.54
1	A	501	THR	CA-CB	-5.03	1.45	1.53
1	B	460	ASP	N-CA	5.02	1.52	1.46
1	B	67	THR	CA-CB	-5.00	1.45	1.53
1	B	77	VAL	CA-C	5.00	1.59	1.52

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	363[A]	HIS	N-CA-C	-8.96	98.45	109.83
1	A	363[B]	HIS	N-CA-C	-8.96	98.45	109.83
1	A	72	LEU	N-CA-C	-8.03	102.61	111.36
1	A	522	ARG	NE-CZ-NH2	7.97	126.37	119.20
1	B	102	SER	CA-C-N	7.69	129.45	119.84
1	B	102	SER	C-N-CA	7.69	129.45	119.84
1	A	578	GLU	N-CA-C	-7.51	97.02	108.96
1	A	268	GLN	N-CA-C	7.41	119.43	111.36
1	A	81	ILE	N-CA-C	7.16	115.26	107.60
1	A	346	HIS	CB-CA-C	7.08	121.86	110.96
1	B	419	GLN	CA-C-O	-6.96	113.98	121.56
1	B	402	GLU	N-CA-C	-6.92	103.73	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	354	ARG	N-CA-C	6.72	118.69	111.36
1	B	268	GLN	N-CA-C	6.70	119.93	111.69
1	B	470	LYS	N-CA-C	-6.69	103.98	111.28
1	A	264	MET	N-CA-C	-6.55	104.22	111.82
1	B	447	PRO	N-CD-CG	-6.55	93.38	103.20
1	B	264	MET	CG-SD-CE	-6.39	86.85	100.90
1	B	44	VAL	CB-CA-C	-6.33	104.45	111.35
1	B	421	ARG	NE-CZ-NH2	6.28	124.85	119.20
1	A	539	PRO	N-CA-C	6.09	121.53	114.03
1	A	40	ARG	CB-CG-CD	5.99	125.07	111.30
1	B	616	GLU	CB-CG-CD	5.97	122.75	112.60
1	B	346	HIS	CB-CA-C	-5.94	101.70	111.68
1	A	266	TRP	N-CA-C	-5.92	104.91	111.36
1	A	410	MET	CB-CA-C	5.91	120.55	110.56
1	A	544	LYS	N-CA-C	-5.89	104.94	111.36
1	A	511	MET	CG-SD-CE	-5.86	88.00	100.90
1	A	601	PRO	N-CA-C	-5.81	103.61	110.70
1	A	466	MET	CG-SD-CE	-5.73	88.29	100.90
1	A	588	GLY	N-CA-C	-5.69	105.49	112.77
1	A	561	ARG	CG-CD-NE	5.60	124.31	112.00
1	B	516	ASN	N-CA-C	-5.58	106.46	113.72
1	B	445	ARG	N-CA-C	-5.55	107.51	114.56
1	B	510	ARG	NE-CZ-NH2	5.55	124.19	119.20
1	B	606	MET	N-CA-C	5.53	117.47	109.07
1	B	89	ASP	N-CA-C	5.52	119.86	113.18
1	B	130	SER	CA-CB-OG	-5.50	100.11	111.10
1	B	363[A]	HIS	N-CA-C	-5.48	102.87	109.83
1	B	363[B]	HIS	N-CA-C	-5.48	102.87	109.83
1	B	583	PHE	N-CA-C	5.47	117.44	108.52
1	B	615	THR	N-CA-C	-5.46	105.33	111.28
1	A	61	SER	CA-C-N	5.46	126.00	119.94
1	A	61	SER	C-N-CA	5.46	126.00	119.94
1	A	225	ILE	N-CA-C	-5.46	106.48	111.67
1	B	460	ASP	N-CA-C	5.45	117.93	111.33
1	B	360	CYS	CA-C-N	5.43	128.30	120.38
1	B	360	CYS	C-N-CA	5.43	128.30	120.38
1	B	452	HIS	N-CA-C	5.43	118.49	108.94
1	B	89	ASP	CA-C-N	-5.42	112.83	122.42
1	B	89	ASP	C-N-CA	-5.42	112.83	122.42
1	B	454	LEU	CA-C-O	-5.34	114.96	121.05
1	A	410	MET	N-CA-C	-5.34	106.36	112.92
1	B	588	GLY	N-CA-C	-5.31	107.19	113.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	502	GLU	N-CA-C	5.31	117.06	111.28
1	B	429	ALA	N-CA-C	-5.26	105.55	111.28
1	A	516	ASN	N-CA-C	-5.24	107.05	113.50
1	A	388	TYR	N-CA-C	-5.24	99.78	108.75
1	A	70	MET	CA-CB-CG	5.22	124.54	114.10
1	A	411	ASN	CB-CA-C	-5.22	104.18	111.80
1	A	118	LYS	CD-CE-NZ	5.20	128.54	111.90
1	A	522	ARG	CG-CD-NE	5.18	123.40	112.00
1	A	347	ARG	CB-CG-CD	5.16	123.17	111.30
1	B	456	GLY	N-CA-C	-5.15	104.12	112.31
1	A	91	ILE	N-CA-C	5.13	116.54	111.67
1	A	289	GLN	N-CA-C	-5.08	105.82	111.36
1	A	411	ASN	N-CA-C	5.06	118.42	111.54
1	B	460	ASP	CB-CA-C	-5.04	102.11	110.68
1	A	466	MET	N-CA-C	-5.04	105.86	111.36
1	B	385	ASP	N-CA-CB	5.02	117.42	110.04
1	B	554	THR	CA-C-N	-5.01	114.10	122.21
1	B	554	THR	C-N-CA	-5.01	114.10	122.21

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	161	PRO	Peptide
1	A	230	ALA	Peptide
1	B	102	SER	Peptide
1	B	180	GLY	Peptide

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	4514	0	4448	91	0
1	B	4718	0	4650	114	0
2	A	27	0	12	0	0
2	B	54	0	23	6	0
3	A	361	0	0	24	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	323	0	0	40	1
All	All	9997	0	9133	206	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 11.

All (206) 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:591:MET:SD	1:B:591:MET:CE	2.02	1.47
1:A:70:MET:CE	1:A:70:MET:SD	2.05	1.45
1:B:70:MET:CE	1:B:70:MET:SD	2.08	1.39
1:B:527:MET:CE	1:B:527:MET:SD	2.09	1.39
1:A:370:MET:CE	1:A:370:MET:SD	2.18	1.31
1:B:538:GLU:HB3	1:B:541:HIS:CD2	1.73	1.24
1:B:538:GLU:HB3	1:B:541:HIS:HD2	0.90	1.06
1:B:484:GLU:HB2	3:B:1205:HOH:O	1.57	1.02
1:B:404:LYS:HE3	3:B:1190:HOH:O	1.65	0.97
1:B:538:GLU:CB	1:B:541:HIS:HD2	1.81	0.94
1:B:531:GLU:HB3	3:B:1132:HOH:O	1.65	0.93
1:B:541:HIS:CE1	1:A:288:LEU:HD21	2.05	0.91
1:A:452:HIS:HD2	1:A:510:ARG:HE	1.16	0.89
1:B:538:GLU:CB	1:B:541:HIS:CD2	2.56	0.89
1:A:47:ARG:HH11	1:A:230:ALA:HB3	1.38	0.87
1:B:578:GLU:O	1:B:579:HIS:ND1	2.09	0.86
1:A:392:GLY:O	1:A:395:GLN:NE2	2.08	0.86
1:A:279:ASN:HB3	3:A:1252:HOH:O	1.74	0.86
1:A:452:HIS:CD2	1:A:510:ARG:HE	1.93	0.85
1:B:246:LEU:HD12	3:B:1185:HOH:O	1.78	0.82
1:B:395:GLN:HG3	3:B:1182:HOH:O	1.81	0.80
1:B:450:LEU:HD21	1:B:452:HIS:ND1	1.98	0.77
1:B:392:GLY:O	1:B:395:GLN:NE2	2.17	0.76
1:B:34:HIS:N	3:B:1150:HOH:O	2.20	0.75
1:B:506:HIS:HE1	3:B:904:HOH:O	1.68	0.74
1:A:388:TYR:HD1	3:A:1261:HOH:O	1.70	0.74
1:B:370:MET:HE1	3:B:1164:HOH:O	1.88	0.73
1:B:241:GLU:N	1:B:241:GLU:CD	2.46	0.73
1:A:214:GLN:HG3	3:A:1240:HOH:O	1.87	0.72
1:B:588:GLY:HA2	1:B:591:MET:HE2	1.73	0.70
1:B:484:GLU:CB	3:B:1205:HOH:O	2.23	0.70
1:B:395:GLN:HB2	3:B:1184:HOH:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:LYS:HZ3	1:A:279:ASN:HD21	1.41	0.68
1:B:485:THR:HG23	3:B:1205:HOH:O	1.94	0.68
1:A:621:LEU:O	1:A:624:ALA:HB3	1.94	0.68
1:B:238:TYR:CE1	3:B:1178:HOH:O	2.47	0.68
1:B:556:GLU:CD	3:B:1201:HOH:O	2.38	0.67
1:A:314:HIS:HB2	3:A:1209:HOH:O	1.94	0.67
1:B:404:LYS:HG2	3:B:1190:HOH:O	1.94	0.67
1:A:47:ARG:NH2	1:A:79:HIS:O	2.27	0.66
1:A:244:LEU:O	1:A:248:LYS:HG3	1.96	0.65
1:A:388:TYR:CD1	3:A:1261:HOH:O	2.48	0.65
1:A:522:ARG:O	1:A:522:ARG:NH1	2.30	0.64
1:A:235:LYS:NZ	1:A:279:ASN:HD21	1.96	0.64
1:B:144:GLN:HG2	3:B:1173:HOH:O	1.99	0.63
1:A:349:GLU:CD	3:A:1227:HOH:O	2.41	0.63
2:B:900:ADP:H5'2	2:B:900:ADP:C8	2.33	0.63
1:A:382:GLN:CD	3:A:1242:HOH:O	2.41	0.63
1:B:241:GLU:CD	1:B:241:GLU:H	2.05	0.62
1:A:241:GLU:H	1:A:241:GLU:CD	2.07	0.62
1:B:185:PHE:HD1	1:B:188:ILE:HD12	1.64	0.62
1:B:399:THR:HG22	3:B:1206:HOH:O	1.99	0.62
1:B:238:TYR:OH	3:B:1178:HOH:O	2.16	0.61
1:A:283:ARG:HD2	3:A:1217:HOH:O	2.00	0.61
1:B:386:ARG:HG2	1:B:388:TYR:CE1	2.35	0.61
1:B:433:GLN:OE1	3:B:1177:HOH:O	2.15	0.61
1:B:363[B]:HIS:HD2	1:B:366:ILE:H	1.50	0.60
1:A:452:HIS:HD2	1:A:510:ARG:NE	1.95	0.60
1:B:96:ASN:HD21	1:B:112:ARG:HD2	1.67	0.60
1:A:241:GLU:HB2	3:A:1208:HOH:O	2.01	0.60
1:B:425:HIS:H	1:B:428:HIS:HD2	1.49	0.60
1:B:506:HIS:O	1:B:510:ARG:HD3	2.01	0.60
1:B:541:HIS:ND1	1:A:288:LEU:HD21	2.16	0.60
1:B:484:GLU:CA	3:B:1205:HOH:O	2.50	0.59
1:B:171:LYS:CE	3:B:1210:HOH:O	2.50	0.59
1:B:450:LEU:HD21	1:B:452:HIS:CE1	2.37	0.59
1:B:184:GLY:HA2	1:B:189:ASP:HB2	1.85	0.58
1:B:241:GLU:N	1:B:241:GLU:OE2	2.36	0.58
1:B:57:LEU:HD23	1:B:156:VAL:HB	1.86	0.57
1:A:235:LYS:NZ	1:A:279:ASN:ND2	2.53	0.57
1:B:317:LYS:O	1:B:321:ASP:HB2	2.05	0.57
1:B:171:LYS:NZ	3:B:1210:HOH:O	2.37	0.56
2:B:900:ADP:C8	2:B:900:ADP:C5'	2.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:GLN:CG	3:A:1240:HOH:O	2.49	0.56
1:B:478:GLU:HG3	1:B:607:ALA:HB1	1.87	0.56
1:A:75:TYR:C	1:A:75:TYR:CD1	2.84	0.56
1:A:476[A]:LEU:HD23	1:A:481:LEU:HB2	1.87	0.56
1:B:96:ASN:ND2	1:B:112:ARG:HD2	2.20	0.56
1:B:569:LYS:HB3	3:B:1204:HOH:O	2.05	0.55
1:A:578:GLU:O	1:A:579:HIS:CB	2.54	0.55
1:B:468:ARG:HD2	3:B:911:HOH:O	2.06	0.55
1:A:579:HIS:O	1:A:580:HIS:C	2.50	0.55
1:B:522:ARG:HB3	3:B:1213:HOH:O	2.07	0.55
1:B:174:TYR:CD2	1:B:188:ILE:HD11	2.42	0.54
1:B:541:HIS:HE1	1:A:288:LEU:HD21	1.63	0.54
1:A:476[A]:LEU:CD2	1:A:481:LEU:HB2	2.37	0.54
1:A:468:ARG:O	1:A:472:HIS:ND1	2.41	0.54
1:B:543:ALA:O	1:B:547:THR:HG23	2.07	0.54
1:A:374:ASP:CG	3:A:1255:HOH:O	2.51	0.54
1:B:174:TYR:HD2	1:B:188:ILE:HD11	1.73	0.53
1:B:410:MET:HE2	1:B:448:VAL:CG2	2.37	0.53
1:B:238:TYR:CZ	3:B:1178:HOH:O	2.58	0.53
1:A:87:ASP:HB2	3:A:1118:HOH:O	2.07	0.53
1:B:410:MET:HE2	1:B:448:VAL:HG21	1.91	0.53
1:B:329:MET:HA	1:B:333:ARG:O	2.08	0.53
1:B:422:ASN:ND2	3:B:992:HOH:O	2.40	0.53
2:B:900:ADP:H5'2	2:B:900:ADP:H8	1.73	0.52
1:A:416[A]:SER:HB2	3:A:1232:HOH:O	2.09	0.52
1:B:45:GLY:N	1:A:44:VAL:O	2.34	0.52
1:A:454:LEU:HD23	1:A:454:LEU:C	2.35	0.52
1:A:578:GLU:O	1:A:579:HIS:ND1	2.43	0.52
1:B:395:GLN:CB	3:B:1184:HOH:O	2.56	0.51
1:A:578:GLU:C	1:A:579:HIS:ND1	2.68	0.51
1:A:395:GLN:H	1:A:395:GLN:CD	2.19	0.51
1:B:347:ARG:N	3:B:1162:HOH:O	2.43	0.51
1:B:345:GLU:HB3	3:B:1168:HOH:O	2.11	0.50
1:B:425:HIS:H	1:B:428:HIS:CD2	2.27	0.50
1:A:578:GLU:O	1:A:579:HIS:CG	2.65	0.50
2:B:900:ADP:C5'	2:B:900:ADP:H8	2.25	0.50
1:B:445:ARG:HD3	3:B:1222:HOH:O	2.11	0.50
1:A:363[B]:HIS:HE1	1:A:497:TYR:O	1.94	0.50
1:B:186:THR:HA	1:B:190:SER:HB2	1.93	0.50
1:A:382:GLN:NE2	3:A:1242:HOH:O	2.45	0.50
1:A:118:LYS:NZ	3:A:1105:HOH:O	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:474:ALA:HB2	3:B:1192:HOH:O	2.12	0.49
1:A:70:MET:CE	1:A:70:MET:CG	2.88	0.49
1:A:452:HIS:HE1	3:A:923:HOH:O	1.96	0.49
1:B:395:GLN:CG	3:B:1184:HOH:O	2.60	0.49
1:B:553:ILE:HD13	3:B:1090:HOH:O	2.13	0.49
1:A:51:ARG:HD2	3:A:1060:HOH:O	2.12	0.49
1:B:507:CYS:HA	1:B:518:TYR:CE1	2.48	0.49
1:A:241:GLU:CB	3:A:1208:HOH:O	2.60	0.49
1:A:388:TYR:HB3	3:A:1177:HOH:O	2.12	0.49
1:A:578:GLU:O	1:A:579:HIS:HB2	2.12	0.49
1:B:588:GLY:HA2	1:B:591:MET:CE	2.41	0.48
1:B:476:LEU:HD23	1:B:481:LEU:HB2	1.95	0.48
1:A:420:LEU:HD13	1:A:453:PRO:HA	1.95	0.48
1:B:92:ARG:HH21	1:B:93:GLN:HG3	1.79	0.48
1:A:239[B]:VAL:HG23	1:A:240:PRO:HD2	1.96	0.48
1:A:416[A]:SER:OG	1:A:439:LEU:HD11	2.14	0.48
1:B:399:THR:CG2	3:B:1206:HOH:O	2.59	0.47
1:A:106:ARG:O	1:A:109:ASN:HB3	2.15	0.47
1:B:476:LEU:HD11	1:B:488:VAL:HG21	1.96	0.47
1:B:567:LYS:HB2	1:B:567:LYS:HE2	1.56	0.47
1:A:340:ASN:N	1:A:341:PRO:CD	2.77	0.47
1:A:341:PRO:HA	1:A:378:GLY:O	2.14	0.47
1:A:385:ASP:CG	3:A:1259:HOH:O	2.56	0.47
1:B:427:GLY:N	1:B:572:MET:HG3	2.30	0.47
1:B:416[A]:SER:OG	1:B:439:LEU:HD11	2.14	0.46
1:B:424:VAL:HA	1:B:428:HIS:HD2	1.79	0.46
1:A:93:GLN:HE21	1:A:93:GLN:HA	1.80	0.46
1:B:238:TYR:HE1	3:B:1178:HOH:O	1.90	0.46
1:A:593:LYS:CG	3:A:1253:HOH:O	2.63	0.46
1:A:340:ASN:N	1:A:341:PRO:HD3	2.30	0.46
1:B:179:ALA:O	1:B:180:GLY:O	2.33	0.46
1:B:313:THR:HA	3:B:1111:HOH:O	2.15	0.46
2:B:900:ADP:H2'	2:B:900:ADP:H5'1	1.76	0.46
1:A:136:THR:O	1:A:137:GLN:C	2.56	0.46
1:A:363[B]:HIS:CD2	1:A:364:PRO:HD2	2.51	0.45
1:B:476:LEU:HD11	1:B:488:VAL:CG2	2.46	0.45
1:B:314:HIS:CD2	1:B:375:TRP:HE1	2.35	0.45
1:B:450:LEU:CD2	1:B:452:HIS:ND1	2.76	0.45
1:A:523:ASP:N	1:A:524:PRO:CD	2.80	0.45
1:A:92:ARG:NH1	3:A:1066:HOH:O	2.49	0.44
1:B:75:TYR:C	1:B:75:TYR:CD1	2.95	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ARG:O	1:B:169:ASP:C	2.60	0.44
1:A:115:GLU:O	1:A:119:LEU:HG	2.17	0.44
1:A:506:HIS:O	1:A:510:ARG:HD3	2.17	0.44
1:B:452:HIS:CD2	1:B:491:PHE:HB2	2.52	0.44
1:A:468:ARG:HB3	1:A:472:HIS:CE1	2.51	0.44
1:B:609:LYS:HD2	3:B:1203:HOH:O	2.17	0.44
1:A:57:LEU:HD23	1:A:156:VAL:HB	1.99	0.44
1:A:453:PRO:HD2	1:A:489:ALA:O	2.16	0.44
1:B:248:LYS:NZ	1:B:385:ASP:OD1	2.46	0.44
1:B:310[A]:LEU:HD13	1:B:377:ILE:HD12	1.99	0.44
1:A:567:LYS:HE2	1:A:584:GLU:HB2	2.00	0.44
1:B:440:LEU:HD23	1:B:440:LEU:HA	1.80	0.44
1:B:163:HIS:CE1	3:B:1166:HOH:O	2.70	0.43
1:A:452:HIS:CD2	1:A:510:ARG:NE	2.73	0.43
1:B:566:ASN:OD1	1:B:566:ASN:C	2.60	0.43
1:A:356:TRP:CD1	1:A:363[B]:HIS:CE1	3.05	0.43
1:B:607:ALA:O	1:B:610:ALA:N	2.52	0.43
1:A:329:MET:HG2	3:A:976:HOH:O	2.19	0.43
1:B:312:ALA:HB2	1:B:377:ILE:HD11	2.01	0.43
1:A:235:LYS:HZ2	1:A:279:ASN:ND2	2.17	0.43
1:B:429:ALA:O	1:B:430:LEU:C	2.57	0.43
1:B:249:THR:O	1:B:250:ASP:C	2.62	0.43
1:B:587:SER:C	1:B:589:THR:N	2.76	0.43
1:A:408:LYS:O	1:A:411:ASN:N	2.41	0.43
1:A:593:LYS:HG3	3:A:1253:HOH:O	2.19	0.42
1:B:562:VAL:HA	2:B:900:ADP:C2	2.53	0.42
1:B:386:ARG:HG3	1:B:387:VAL:N	2.34	0.42
1:B:166:GLU:O	1:B:169:ASP:HB2	2.20	0.42
1:A:151:LEU:HA	1:A:152:PRO:HD3	1.92	0.42
1:A:425:HIS:O	1:A:428:HIS:HB2	2.19	0.42
1:B:233:GLU:O	1:B:234:VAL:C	2.61	0.42
1:A:529:HIS:HA	1:A:530:PRO:HD3	1.93	0.42
1:B:239:VAL:HG12	1:B:383:VAL:O	2.20	0.41
1:B:112:ARG:O	1:B:116:VAL:HG23	2.21	0.41
1:B:538:GLU:OE2	1:B:540:SER:OG	2.37	0.41
1:A:511:MET:SD	1:A:555[A]:LEU:HD22	2.61	0.41
1:B:244:LEU:O	1:B:248:LYS:HG3	2.20	0.41
1:B:484:GLU:C	3:B:1205:HOH:O	2.62	0.41
1:B:471:GLN:O	1:B:475:VAL:HG23	2.21	0.41
1:A:623:LYS:O	1:A:624:ALA:C	2.64	0.41
1:B:171:LYS:HE2	3:B:1210:HOH:O	2.16	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:VAL:HA	1:A:227:PRO:HD3	1.91	0.41
1:B:151:LEU:HA	1:B:152:PRO:HD3	1.90	0.41
1:A:356:TRP:NE1	1:A:363[A]:HIS:CE1	2.89	0.40
1:A:347:ARG:HH11	1:A:347:ARG:HD2	1.73	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)
3:B:1218:HOH:O	3:A:1007:HOH:O[1_455]	2.00	0.20

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	566/630 (90%)	551 (97%)	14 (2%)	1 (0%)	43	38
1	B	592/630 (94%)	567 (96%)	19 (3%)	6 (1%)	12	5
All	All	1158/1260 (92%)	1118 (96%)	33 (3%)	7 (1%)	21	13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	180	GLY
1	B	183	LYS
1	B	579	HIS
1	A	579	HIS
1	B	597	GLU
1	B	169	ASP
1	B	181	GLU

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	489/538 (91%)	454 (93%)	35 (7%)	13	7
1	B	508/538 (94%)	470 (92%)	38 (8%)	12	6
All	All	997/1076 (93%)	924 (93%)	73 (7%)	13	6

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	38	ASN
1	B	106	ARG
1	B	144	GLN
1	B	167	GLN
1	B	173	LEU
1	B	175	LYS
1	B	182	ILE
1	B	183	LYS
1	B	186	THR
1	B	241	GLU
1	B	246	LEU
1	B	277	PRO
1	B	304	LEU
1	B	310[A]	LEU
1	B	310[B]	LEU
1	B	315	GLU
1	B	323	CYS
1	B	361	LYS
1	B	374	ASP
1	B	408	LYS
1	B	420	LEU
1	B	437	LYS
1	B	445	ARG
1	B	508	ARG
1	B	522	ARG
1	B	531	GLU
1	B	555	LEU

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Mol	Chain	Res	Type
1	B	562	VAL
1	B	567	LYS
1	B	569	LYS
1	B	570	LYS
1	B	578	GLU
1	B	581	GLU
1	B	586	ILE
1	B	593	LYS
1	B	621	LEU
1	B	622	GLU
1	B	623	LYS
1	A	39	LYS
1	A	42	GLN
1	A	44	VAL
1	A	46	THR
1	A	70	MET
1	A	93	GLN
1	A	214	GLN
1	A	222	GLU
1	A	241	GLU
1	A	246[A]	LEU
1	A	246[B]	LEU
1	A	248	LYS
1	A	304	LEU
1	A	347	ARG
1	A	361	LYS
1	A	395	GLN
1	A	408	LYS
1	A	410	MET
1	A	420	LEU
1	A	441	GLU
1	A	445	ARG
1	A	452	HIS
1	A	472	HIS
1	A	508	ARG
1	A	522	ARG
1	A	553	ILE
1	A	555[A]	LEU
1	A	555[B]	LEU
1	A	570	LYS
1	A	577	SER
1	A	578	GLU

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Mol	Chain	Res	Type
1	A	581	GLU
1	A	596	ARG
1	A	603	GLU
1	A	608	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	38	ASN
1	B	96	ASN
1	B	109	ASN
1	B	144	GLN
1	B	210	ASN
1	B	314	HIS
1	B	411	ASN
1	B	422	ASN
1	B	428	HIS
1	B	438	GLN
1	B	506	HIS
1	B	541	HIS
1	B	599	GLN
1	A	42	GLN
1	A	93	GLN
1	A	140	ASN
1	A	144	GLN
1	A	279	ASN
1	A	372	GLN
1	A	422	ASN
1	A	436	HIS
1	A	452	HIS
1	A	580	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	A	900	-	28,29,29	1.50	5 (17%)	43,45,45	1.95	9 (20%)
2	ADP	B	800	-	28,29,29	1.56	7 (25%)	43,45,45	1.46	9 (20%)
2	ADP	B	900	-	28,29,29	1.85	8 (28%)	43,45,45	3.29	22 (51%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	A	900	-	-	5/16/32/32	0/3/3/3
2	ADP	B	800	-	-	4/16/32/32	0/3/3/3
2	ADP	B	900	-	-	7/16/32/32	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	ADP	O5'-C5'	-4.45	1.27	1.44
2	A	900	ADP	PA-O3A	-4.02	1.55	1.59
2	B	800	ADP	C2-N3	3.85	1.40	1.33
2	B	900	ADP	PB-O3B	-3.84	1.40	1.54
2	B	800	ADP	PA-O3A	3.34	1.63	1.59
2	B	900	ADP	C2-N1	3.27	1.39	1.33
2	B	900	ADP	O4'-C4'	-2.88	1.38	1.45
2	B	800	ADP	O2'-C2'	-2.86	1.35	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	800	ADP	C2-N1	2.73	1.38	1.33
2	A	900	ADP	C4-N9	-2.58	1.32	1.37
2	A	900	ADP	C2-N1	2.50	1.38	1.33
2	B	900	ADP	C5-N7	-2.47	1.34	1.39
2	B	800	ADP	O4'-C4'	-2.46	1.39	1.45
2	A	900	ADP	O2'-C2'	-2.45	1.36	1.43
2	B	900	ADP	C5-C6	-2.31	1.34	1.41
2	B	900	ADP	PA-O5'	-2.21	1.50	1.59
2	B	800	ADP	C4-N3	2.16	1.38	1.34
2	B	800	ADP	PA-O2A	-2.03	1.45	1.55
2	B	900	ADP	C8-N7	2.03	1.35	1.31
2	A	900	ADP	O5'-C5'	-2.03	1.36	1.44

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	900	ADP	O2A-PA-O5'	8.35	145.42	107.57
2	B	900	ADP	O3A-PA-O1A	-7.70	87.53	110.70
2	B	900	ADP	O4'-C4'-C5'	-7.10	86.60	109.33
2	B	900	ADP	N3-C2-N1	-5.60	120.10	128.58
2	A	900	ADP	N3-C2-N1	-5.03	120.97	128.58
2	B	900	ADP	N3-C4-N9	4.90	135.51	127.17
2	B	900	ADP	C5-C4-N3	-4.89	119.98	126.72
2	A	900	ADP	O3B-PB-O2B	4.85	126.00	107.80
2	B	900	ADP	O2A-PA-O3A	-4.60	94.84	107.27
2	B	900	ADP	C5-C6-N6	-4.49	112.18	123.29
2	B	900	ADP	O3B-PB-O2B	4.33	124.05	107.80
2	A	900	ADP	N9-C8-N7	-4.02	108.23	113.94
2	B	900	ADP	N9-C8-N7	-4.00	108.25	113.94
2	A	900	ADP	C4-N9-C8	3.95	109.89	105.74
2	A	900	ADP	O2A-PA-O3A	3.79	117.51	107.27
2	B	900	ADP	O2B-PB-O1B	-3.65	96.61	110.83
2	A	900	ADP	C2-N1-C6	3.57	124.59	118.73
2	B	900	ADP	N6-C6-N1	3.46	126.08	118.38
2	B	900	ADP	C2-N3-C4	3.43	120.20	111.83
2	B	800	ADP	N3-C2-N1	-3.27	123.63	128.58
2	B	900	ADP	O2'-C2'-C3'	3.22	122.15	111.82
2	B	900	ADP	C2'-C1'-N9	3.17	121.19	113.30
2	A	900	ADP	O4'-C1'-N9	-3.13	102.07	108.09
2	B	900	ADP	C6-C5-C4	2.87	121.09	117.18
2	B	900	ADP	O5'-PA-O1A	-2.84	97.67	108.94
2	B	900	ADP	C6-C5-N7	-2.81	126.68	132.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	800	ADP	N9-C8-N7	-2.80	109.96	113.94
2	B	800	ADP	O2B-PB-O1B	2.77	121.61	110.83
2	B	800	ADP	C4-N9-C8	2.62	108.49	105.74
2	B	900	ADP	O3B-PB-O1B	2.49	120.53	110.83
2	B	800	ADP	O2A-PA-O5'	2.45	118.65	107.57
2	B	800	ADP	O2A-PA-O3A	2.44	113.87	107.27
2	A	900	ADP	O3A-PB-O1B	-2.40	98.41	111.04
2	B	900	ADP	C5-N7-C8	2.28	107.03	103.45
2	B	800	ADP	O4'-C1'-N9	2.21	112.33	108.09
2	B	800	ADP	C5-N7-C8	2.21	106.92	103.45
2	B	900	ADP	C4-N9-C8	2.20	108.05	105.74
2	B	900	ADP	O4'-C1'-N9	-2.15	103.95	108.09
2	A	900	ADP	O3A-PA-O1A	-2.05	104.53	110.70
2	B	800	ADP	O3A-PB-O1B	-2.00	100.50	111.04

There are no chirality outliers.

All (16) torsion outliers are listed below:

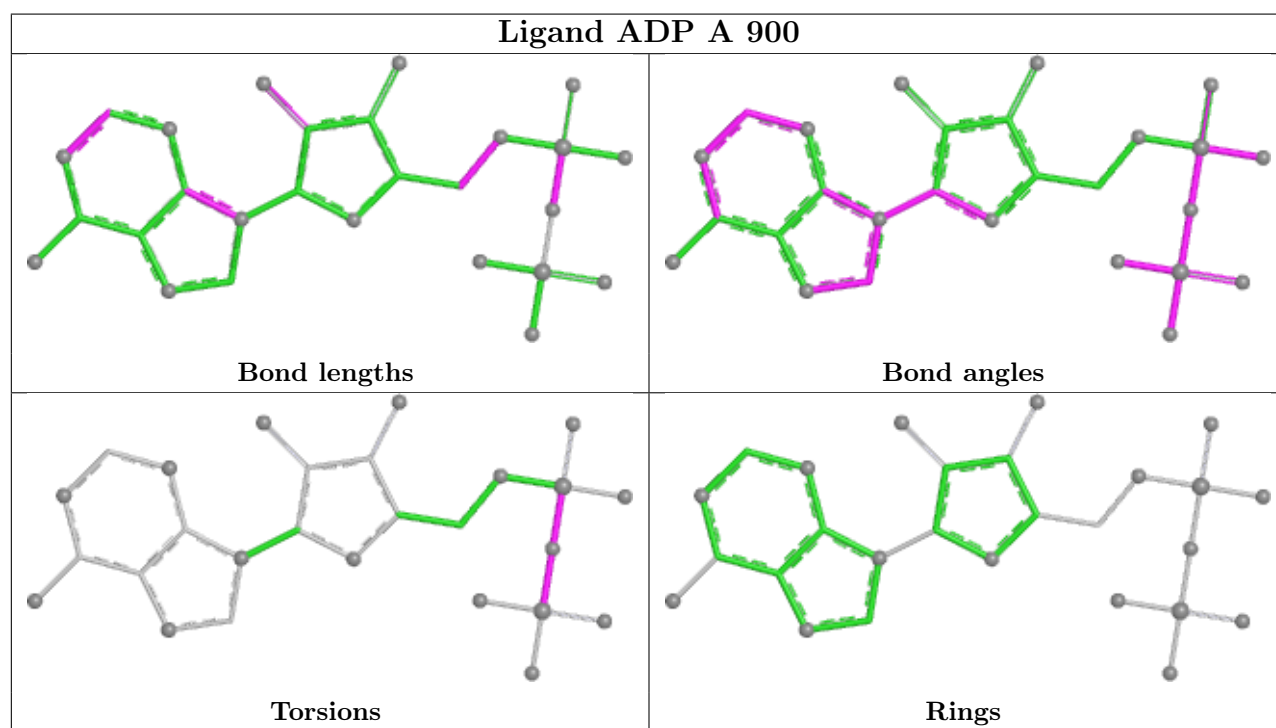
Mol	Chain	Res	Type	Atoms
2	B	800	ADP	PA-O3A-PB-O2B
2	B	900	ADP	PA-O3A-PB-O3B
2	B	900	ADP	C5'-O5'-PA-O2A
2	B	900	ADP	C5'-O5'-PA-O3A
2	A	900	ADP	PA-O3A-PB-O3B
2	B	900	ADP	C3'-C4'-C5'-O5'
2	B	900	ADP	O4'-C4'-C5'-O5'
2	B	800	ADP	PA-O3A-PB-O1B
2	A	900	ADP	PB-O3A-PA-O5'
2	B	900	ADP	C5'-O5'-PA-O1A
2	A	900	ADP	PB-O3A-PA-O1A
2	A	900	ADP	PA-O3A-PB-O1B
2	A	900	ADP	PA-O3A-PB-O2B
2	B	800	ADP	O4'-C4'-C5'-O5'
2	B	800	ADP	O4'-C1'-N9-C8
2	B	900	ADP	PA-O3A-PB-O1B

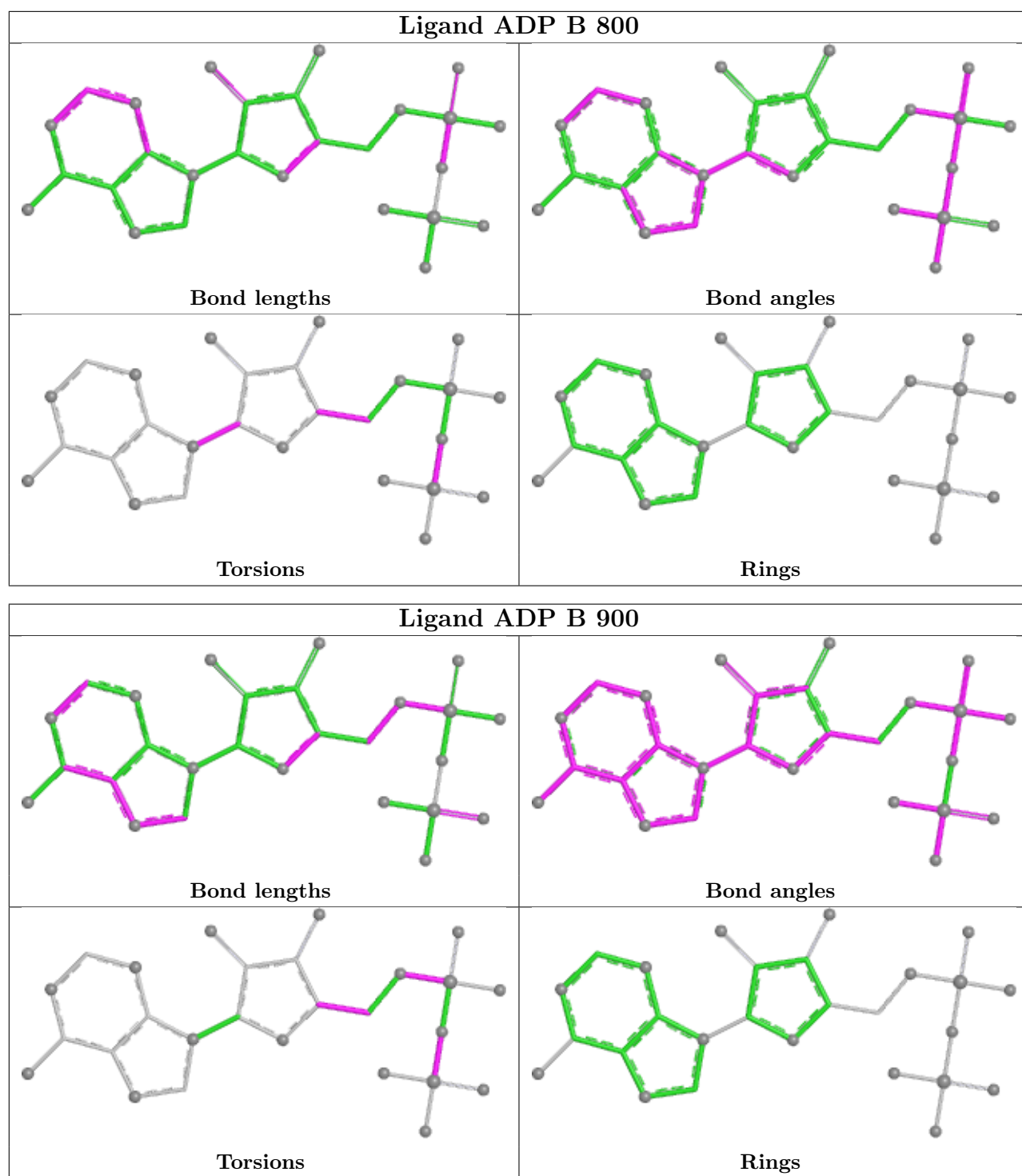
There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	900	ADP	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	563/630 (89%)	-0.19	8 (1%) 73 75	16, 37, 67, 90	10 (1%)
1	B	590/630 (93%)	0.00	9 (1%) 72 74	20, 39, 72, 104	4 (0%)
All	All	1153/1260 (91%)	-0.09	17 (1%) 72 74	16, 38, 69, 104	14 (1%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	TYR	4.5
1	A	162	LEU	4.4
1	B	180	GLY	3.0
1	A	46	THR	2.8
1	B	38	ASN	2.6
1	B	388	TYR	2.6
1	B	182	ILE	2.5
1	A	579	HIS	2.5
1	A	596	ARG	2.5
1	A	230	ALA	2.3
1	B	179	ALA	2.2
1	A	246[A]	LEU	2.2
1	B	245	HIS	2.1
1	B	185	PHE	2.1
1	B	41	GLY	2.1
1	A	34	HIS	2.0
1	B	184	GLY	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](#)

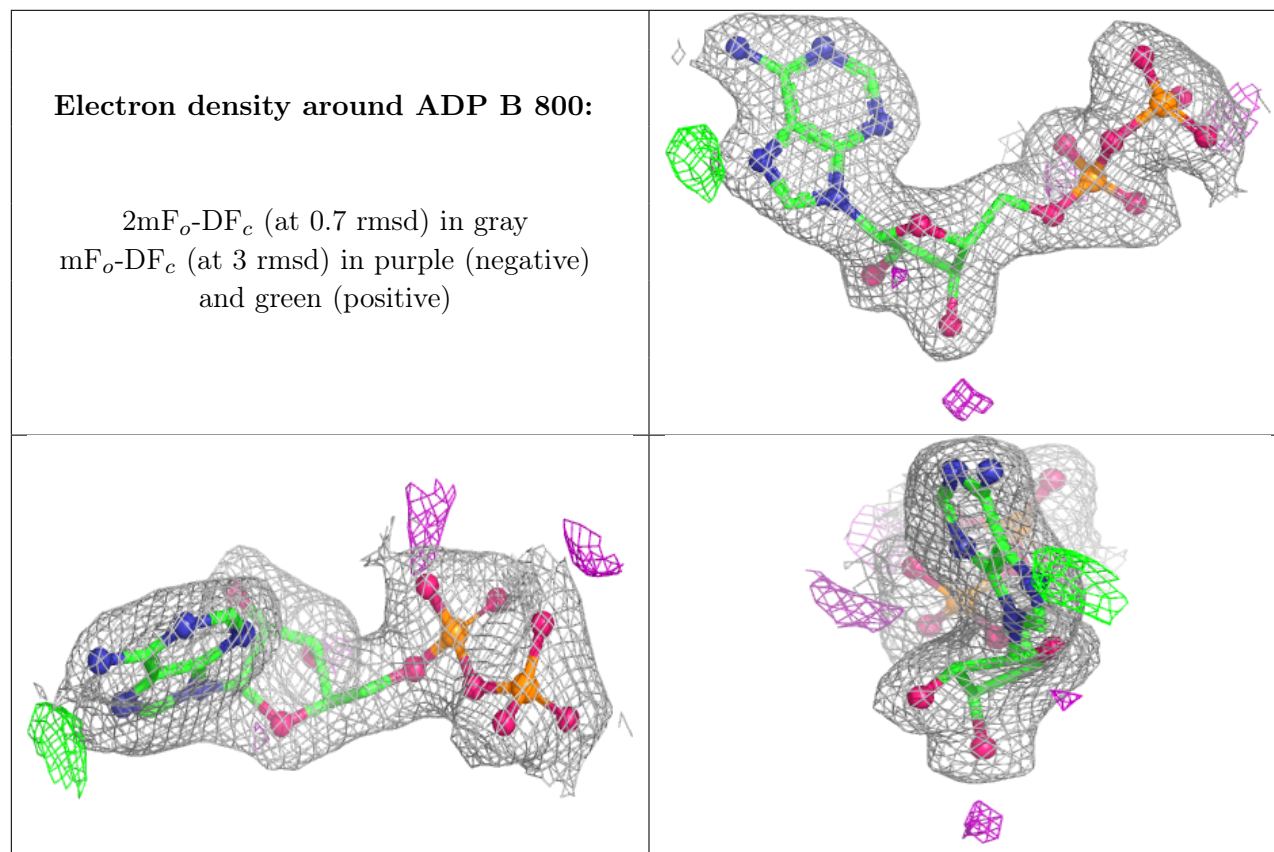
There are no oligosaccharides in this entry.

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.

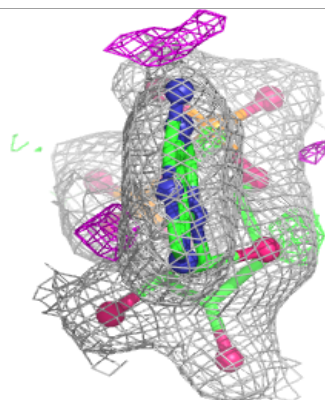
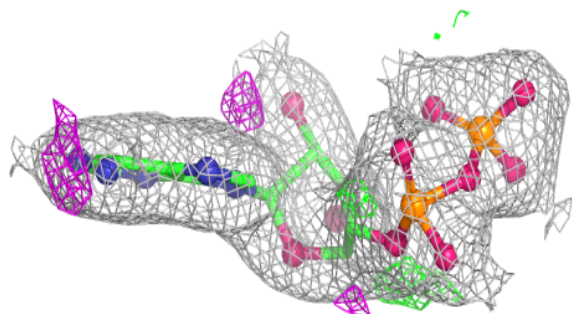
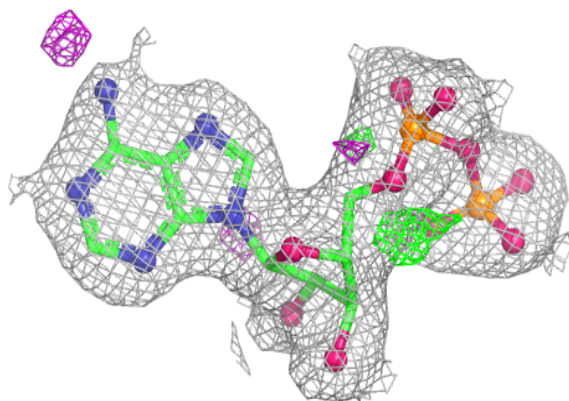
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ADP	B	800	27/27	0.98	0.05	37,41,43,49	0
2	ADP	B	900	27/27	0.98	0.06	33,38,46,48	0
2	ADP	A	900	27/27	0.99	0.04	30,33,39,41	0

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.

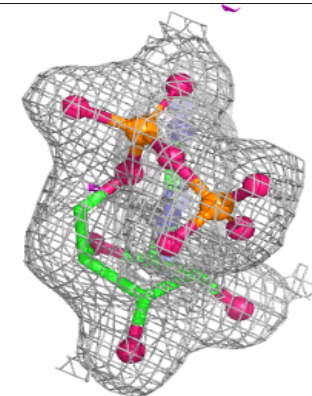
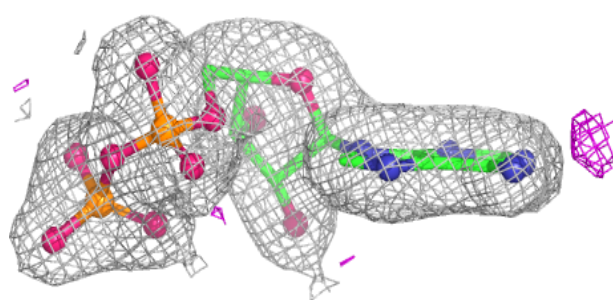
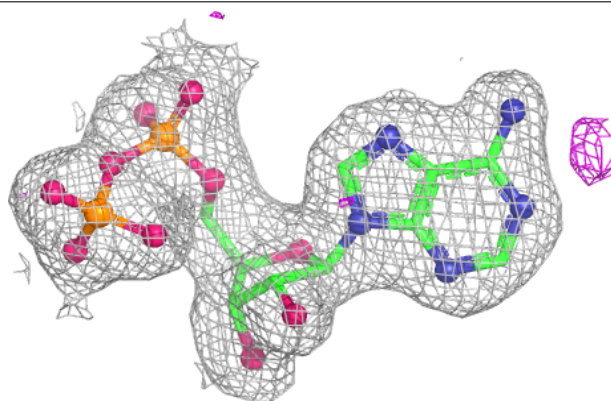


Electron density around ADP B 900:

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

**Electron density around ADP A 900:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.