



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:18 AM GMT

PDB ID : 3LL5
Title : Crystal structure of T. acidophilum isopentenyl phosphate kinase product complex
Authors : Mabanglo, M.F.; Hill, C.P.
Deposited on : 2010-01-28
Resolution : 1.99 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

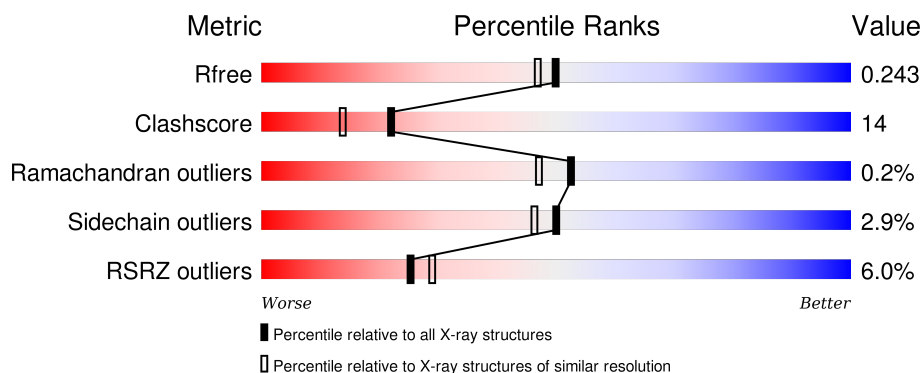
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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}	91344	8664 (2.00-1.96)
Clashscore	102246	9905 (2.00-1.96)
Ramachandran outliers	100387	9792 (2.00-1.96)
Sidechain outliers	100360	9791 (2.00-1.96)
RSRZ outliers	91569	8679 (2.00-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	249	<div> <div>2%</div> <div>76% 16% • 6%</div> </div>
1	B	249	<div> <div>3%</div> <div>78% 16% 6%</div> </div>
1	C	249	<div> <div>4%</div> <div>73% 20% • 5%</div> </div>
1	D	249	<div> <div>12%</div> <div>69% 24% • 6%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	IP8	C	247	-	-	X	X
5	IP8	D	247	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7677 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 Gamma-glutamyl kinase related protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	Se	2	0	0
			1804	1149	304	338	1	12			
1	B	234	Total	C	N	O	S	Se	1	0	0
			1808	1151	305	339	1	12			
1	C	237	Total	C	N	O	S	Se	1	1	0
			1839	1173	308	344	1	13			
1	D	234	Total	C	N	O	S	Se	1	1	0
			1822	1163	307	339	1	12			

There are 16 discrepancies between the modelled and reference sequences:

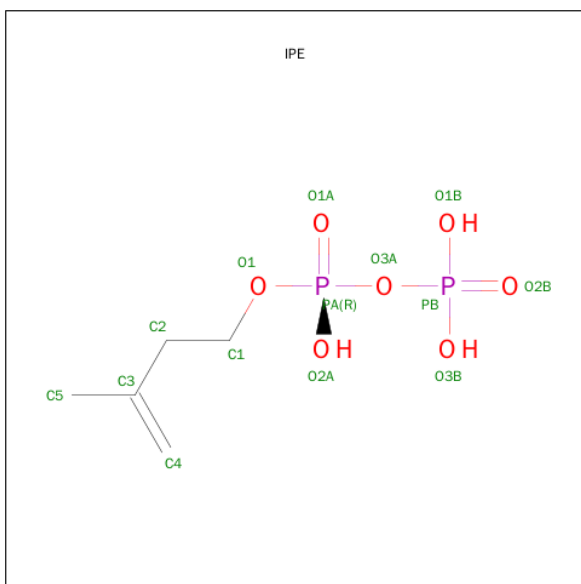
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ASP	-	EXPRESSION TAG	UNP Q9HLX1
A	-2	PRO	-	EXPRESSION TAG	UNP Q9HLX1
A	-1	PHE	-	EXPRESSION TAG	UNP Q9HLX1
A	0	THR	-	EXPRESSION TAG	UNP Q9HLX1
B	-3	ASP	-	EXPRESSION TAG	UNP Q9HLX1
B	-2	PRO	-	EXPRESSION TAG	UNP Q9HLX1
B	-1	PHE	-	EXPRESSION TAG	UNP Q9HLX1
B	0	THR	-	EXPRESSION TAG	UNP Q9HLX1
C	-3	ASP	-	EXPRESSION TAG	UNP Q9HLX1
C	-2	PRO	-	EXPRESSION TAG	UNP Q9HLX1
C	-1	PHE	-	EXPRESSION TAG	UNP Q9HLX1
C	0	THR	-	EXPRESSION TAG	UNP Q9HLX1
D	-3	ASP	-	EXPRESSION TAG	UNP Q9HLX1
D	-2	PRO	-	EXPRESSION TAG	UNP Q9HLX1
D	-1	PHE	-	EXPRESSION TAG	UNP Q9HLX1
D	0	THR	-	EXPRESSION TAG	UNP Q9HLX1

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	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

- Molecule 3 is 3-METHYLBUT-3-ENYL TRIHYDROGEN DIPHOSPHATE (three-letter code: IPE) (formula: $\text{C}_5\text{H}_{12}\text{O}_7\text{P}_2$).



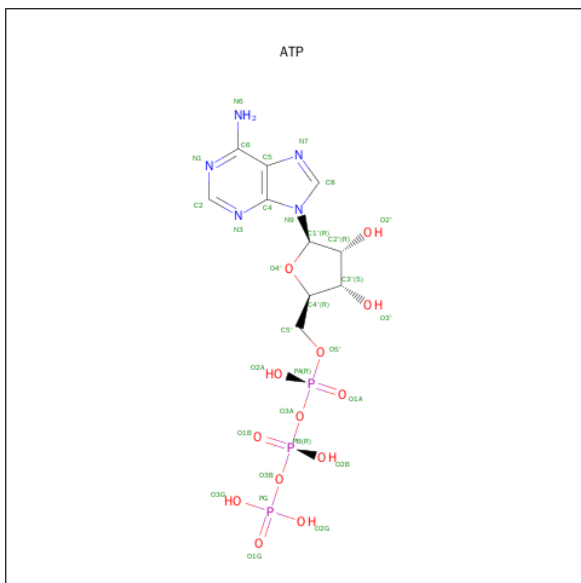
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			14	5	7	2		

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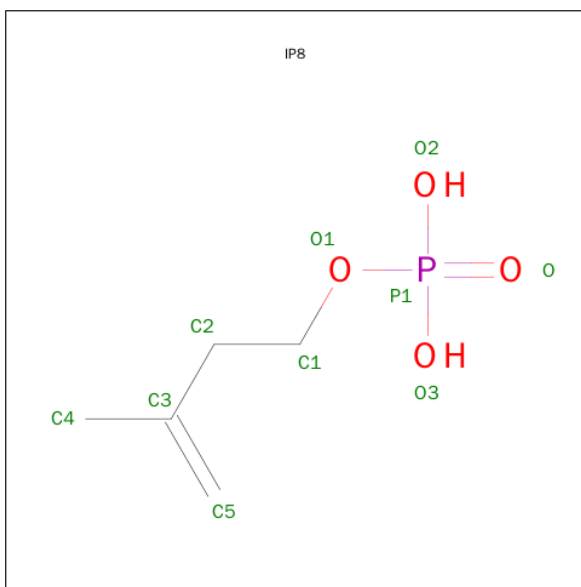
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O	P	0	0
			14	5	7	2		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0
			31	10	5	13	3	
4	D	1	Total	C	N	O	P	0
			31	10	5	13	3	

- Molecule 5 is ISOPENTENYL PHOSPHATE (three-letter code: IP8) (formula: $C_5H_{11}O_4P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	O	P	0	0
			10	5	4	1		
5	D	1	Total	C	O	P	0	0
			10	5	4	1		

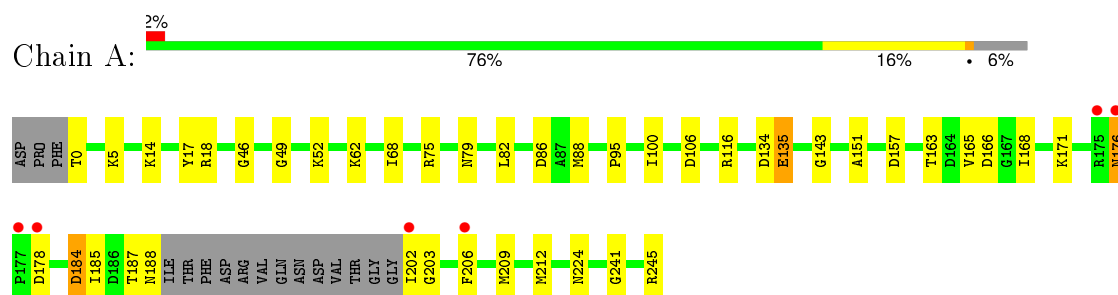
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	65	Total	O	0	0
			65	65		
6	B	58	Total	O	0	0
			58	58		
6	C	70	Total	O	0	0
			70	70		
6	D	47	Total	O	0	0
			47	47		

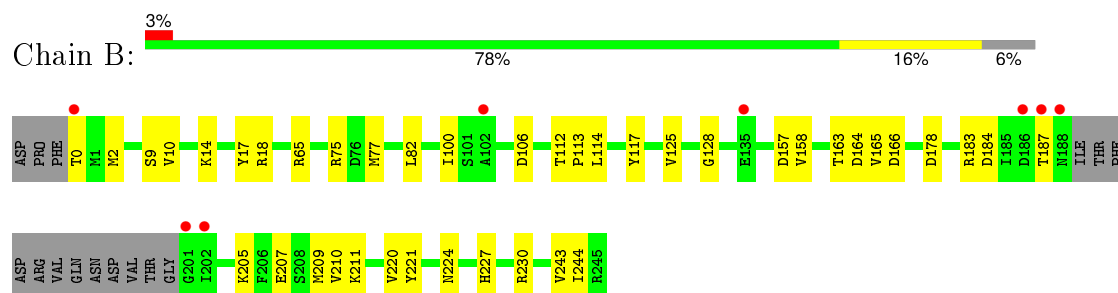
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

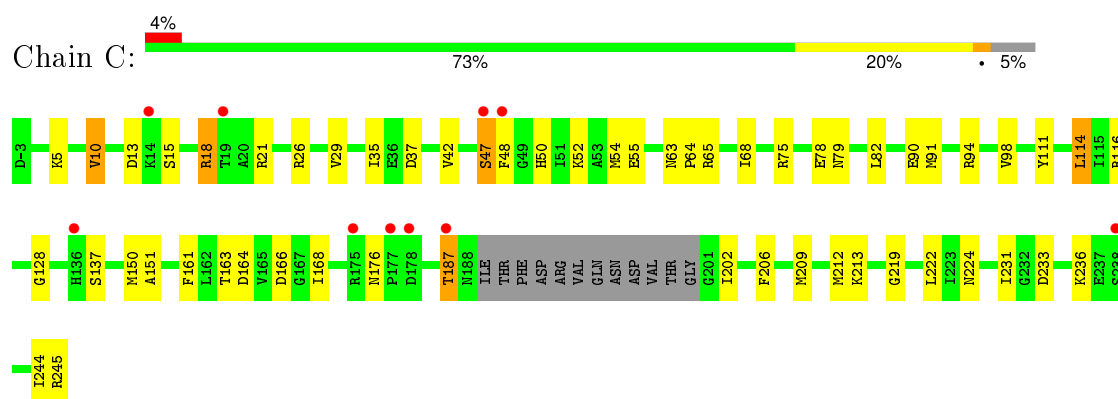
- Molecule 1: Gamma-glutamyl kinase related protein



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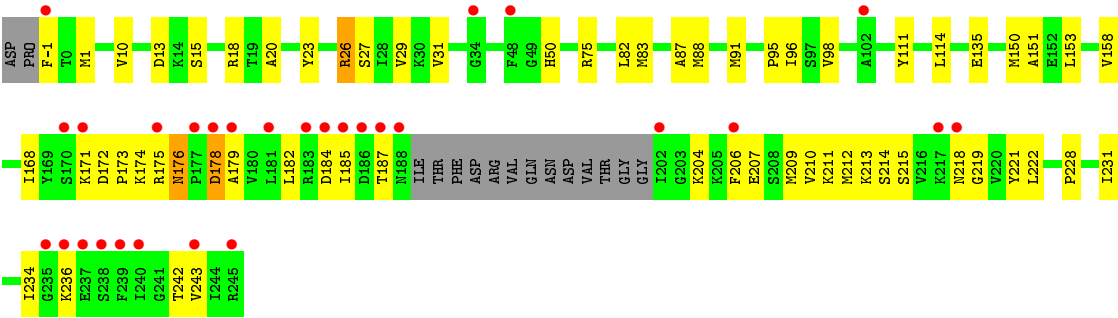


- Molecule 1: Gamma-glutamyl kinase related protein



- Molecule 1: Gamma-glutamyl kinase related protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	187.10 Å 42.79 Å 134.90 Å 90.00° 113.19° 90.00°	Depositor
Resolution (Å)	38.29 – 1.99 38.29 – 1.99	Depositor EDS
% Data completeness (in resolution range)	96.0 (38.29-1.99) 96.0 (38.29-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 1.98 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.193 , 0.238 0.198 , 0.243	Depositor DCC
R_{free} test set	3342 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 65832 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7677	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: IP8, ATP, ADP, IPE

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	0.98	1/1825 (0.1%)	0.91	3/2440 (0.1%)
1	B	0.96	2/1829 (0.1%)	0.91	1/2445 (0.0%)
1	C	1.01	1/1865 (0.1%)	0.96	5/2494 (0.2%)
1	D	0.96	1/1848 (0.1%)	0.91	2/2471 (0.1%)
All	All	0.98	5/7367 (0.1%)	0.92	11/9850 (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	C	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	106	ASP	CB-CG	8.25	1.69	1.51
1	B	125	VAL	CB-CG2	5.64	1.64	1.52
1	B	117	TYR	CE1-CZ	5.44	1.45	1.38
1	D	135	GLU	CG-CD	5.13	1.59	1.51
1	C	10	VAL	CB-CG1	-5.03	1.42	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	21	ARG	NE-CZ-NH2	-10.07	115.26	120.30
1	A	116	ARG	NE-CZ-NH2	-7.40	116.60	120.30
1	D	75	ARG	NE-CZ-NH2	6.96	123.78	120.30
1	B	65	ARG	NE-CZ-NH1	6.11	123.35	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	21	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	A	116	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	C	37	ASP	CB-CG-OD1	5.56	123.30	118.30
1	C	116	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	106	ASP	CB-CG-OD1	5.31	123.08	118.30
1	C	48	PHE	N-CA-CB	5.25	120.05	110.60
1	D	184	ASP	C-N-CA	-5.23	108.62	121.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	47	SER	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	1804	0	1831	40	0
1	B	1808	0	1834	32	0
1	C	1839	0	1863	57	0
1	D	1822	0	1847	58	0
2	A	27	0	12	4	0
2	B	27	0	12	3	0
3	A	14	0	9	5	0
3	B	14	0	9	2	0
4	C	31	0	12	6	0
4	D	31	0	12	5	0
5	C	10	0	9	13	0
5	D	10	0	9	11	0
6	A	65	0	0	6	0
6	B	58	0	0	0	0
6	C	70	0	0	11	0
6	D	47	0	0	9	0
All	All	7677	0	7459	207	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 14.

All (207) 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:D:83:MSE:CE	1:D:83:MSE:SE	2.14	1.46
1:D:1:MSE:CE	1:D:1:MSE:SE	2.15	1.45
4:D:246:ATP:O1G	5:D:247:IP8:P1	2.07	1.11
4:C:246:ATP:O1G	5:C:247:IP8:P1	2.09	1.10
3:A:247:IPE:H11	6:A:291:HOH:O	1.53	1.07
1:D:182:LEU:CD1	1:D:185:ILE:HD11	1.88	1.04
1:C:50:HIS:HB3	6:C:296:HOH:O	1.62	0.97
5:D:247:IP8:H5A	6:D:263:HOH:O	1.67	0.95
1:D:182:LEU:HD13	1:D:185:ILE:HD11	1.44	0.95
5:D:247:IP8:C5	6:D:263:HOH:O	2.20	0.89
1:D:231:ILE:O	1:D:234:ILE:HD12	1.73	0.89
1:C:98:VAL:HB	1:C:114:LEU:HD21	1.52	0.89
2:A:246:ADP:O3B	3:A:247:IPE:PB	2.31	0.88
5:D:247:IP8:H1	6:D:272:HOH:O	1.76	0.84
1:C:209:MSE:HE3	1:C:244:ILE:CD1	2.10	0.81
1:C:233:ASP:HA	1:C:236:LYS:HD3	1.62	0.80
1:A:18:ARG:NH2	1:A:79:ASN:OD1	2.15	0.79
2:B:246:ADP:O3B	3:B:247:IPE:PB	2.41	0.77
5:C:247:IP8:C1	6:C:271:HOH:O	2.31	0.77
1:C:26:ARG:NH2	1:C:90:GLU:OE2	2.17	0.77
1:D:182:LEU:HD12	1:D:185:ILE:HD11	1.64	0.77
1:D:206:PHE:HA	1:D:209:MSE:CE	2.15	0.77
4:D:246:ATP:O1G	5:D:247:IP8:O3	2.03	0.76
1:A:171:LYS:HB2	1:A:176:ASN:OD1	1.86	0.76
4:D:246:ATP:PG	5:D:247:IP8:O3	2.44	0.75
5:C:247:IP8:H1A	6:C:271:HOH:O	1.87	0.74
1:C:50:HIS:HD2	5:C:247:IP8:H4A	1.51	0.73
1:D:26:ARG:HH11	1:D:26:ARG:HG2	1.54	0.73
1:C:50:HIS:CD2	5:C:247:IP8:H4A	2.23	0.72
1:C:111:TYR:CG	1:C:150:MSE:HE1	2.26	0.71
1:C:213:LYS:HE3	1:C:245:ARG:HA	1.73	0.71
1:C:213:LYS:HE2	1:C:219:GLY:HA2	1.72	0.70
1:D:206:PHE:HA	1:D:209:MSE:HE3	1.72	0.70
1:B:163:THR:HG23	1:B:165:VAL:H	1.57	0.69
1:D:221:TYR:CE2	1:D:243:VAL:HG22	2.28	0.69
1:A:171:LYS:CB	1:A:176:ASN:OD1	2.40	0.68
1:C:10:VAL:CG1	1:C:163:THR:C	2.62	0.68
1:D:182:LEU:HD13	1:D:185:ILE:CD1	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:221:TYR:CD2	1:D:243:VAL:HG22	2.29	0.67
1:C:26:ARG:CZ	1:C:90:GLU:OE1	2.44	0.66
1:A:168:ILE:HD11	1:A:209:MSE:HE1	1.78	0.66
1:C:111:TYR:CD2	1:C:150:MSE:HE1	2.30	0.65
1:D:213:LYS:CD	1:D:219:GLY:HA2	2.27	0.65
1:D:111:TYR:CG	1:D:150:MSE:HE1	2.31	0.65
1:C:18:ARG:HD2	6:C:309:HOH:O	1.95	0.65
1:A:202:ILE:N	6:A:289:HOH:O	2.29	0.65
1:D:213:LYS:HD3	1:D:219:GLY:HA2	1.78	0.65
4:C:246:ATP:PG	5:C:247:IP8:O3	2.55	0.65
1:C:98:VAL:CB	1:C:114:LEU:HD21	2.24	0.65
5:C:247:IP8:H1	6:C:271:HOH:O	1.96	0.65
5:D:247:IP8:H4A	6:D:293:HOH:O	1.98	0.64
1:C:128:GLY:HA3	6:C:279:HOH:O	1.98	0.64
5:D:247:IP8:C1	6:D:272:HOH:O	2.36	0.64
1:C:187:THR:HG21	1:C:213:LYS:HG2	1.80	0.64
5:C:247:IP8:H4B	6:C:275:HOH:O	1.97	0.63
1:A:168:ILE:HD13	1:A:202:ILE:HD11	1.81	0.63
1:D:111:TYR:CD2	1:D:150:MSE:HE1	2.33	0.63
2:B:246:ADP:O3B	3:B:247:IPE:O2B	2.17	0.62
1:A:184:ASP:OD1	1:A:245:ARG:HD2	1.99	0.62
1:B:9:SER:HB3	1:B:14:LYS:HE3	1.81	0.62
1:C:187:THR:HG22	1:C:244:ILE:HG22	1.81	0.62
1:C:187:THR:HG23	1:C:245:ARG:C	2.21	0.60
1:A:14:LYS:HD3	6:A:287:HOH:O	2.00	0.59
4:D:246:ATP:O1G	5:D:247:IP8:O2	2.20	0.59
1:C:18:ARG:CD	6:C:309:HOH:O	2.50	0.59
1:C:52:LYS:O	1:C:55:GLU:HG2	2.02	0.59
1:D:151:ALA:HB2	1:D:212:MSE:HE3	1.85	0.58
1:A:5:LYS:HE3	1:A:143:GLY:HA3	1.85	0.58
1:C:78:GLU:HG2	6:C:299:HOH:O	2.03	0.58
1:C:90:GLU:C	1:C:91:MSE:HE2	2.25	0.57
1:C:18:ARG:O	6:C:309:HOH:O	2.18	0.57
1:D:26:ARG:HH11	1:D:26:ARG:CG	2.17	0.57
1:B:106:ASP:C	1:B:106:ASP:OD1	2.43	0.57
1:C:10:VAL:HG11	1:C:163:THR:C	2.26	0.56
1:A:202:ILE:HG21	2:A:246:ADP:N7	2.21	0.56
1:B:82:LEU:CD1	1:D:82:LEU:HD12	2.36	0.56
1:B:10:VAL:CG2	1:B:163:THR:C	2.74	0.55
1:C:65:ARG:HD2	6:C:288:HOH:O	2.06	0.55
1:A:168:ILE:CD1	1:A:202:ILE:HD11	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:0:THR:HA	1:A:157:ASP:OD1	2.07	0.55
1:C:206:PHE:HA	1:C:209:MSE:HE2	1.87	0.55
1:C:168:ILE:CD1	1:C:202:ILE:HD13	2.37	0.55
4:C:246:ATP:O1G	5:C:247:IP8:O3	2.25	0.55
1:D:29:VAL:HG12	1:D:91:MSE:HG3	1.88	0.54
1:D:50[B]:HIS:CD2	6:D:293:HOH:O	2.61	0.54
1:D:207:GLU:O	1:D:211:LYS:HG3	2.07	0.54
1:A:163:THR:HG23	1:A:165:VAL:H	1.72	0.54
1:B:209:MSE:HE2	1:B:244:ILE:HD13	1.90	0.54
1:B:209:MSE:HE3	1:B:220:VAL:HG12	1.90	0.54
1:D:206:PHE:HA	1:D:209:MSE:HE2	1.89	0.54
1:B:82:LEU:HD12	1:D:82:LEU:CD1	2.38	0.53
1:D:171:LYS:HD2	1:D:176:ASN:ND2	2.24	0.53
1:A:75:ARG:HD2	1:A:75:ARG:C	2.29	0.53
1:B:17:TYR:CZ	1:B:18:ARG:HD2	2.44	0.53
1:A:62:LYS:HD2	1:A:135:GLU:O	2.08	0.53
1:A:151:ALA:HB2	1:A:212:MSE:HE3	1.91	0.53
1:C:50:HIS:HE2	5:C:247:IP8:P1	2.32	0.52
1:B:82:LEU:CD1	1:D:82:LEU:CD1	2.87	0.52
1:C:111:TYR:CE1	1:C:150:MSE:HE3	2.46	0.51
1:A:0:THR:N	6:A:295:HOH:O	2.42	0.51
4:C:246:ATP:O1G	5:C:247:IP8:O	2.29	0.51
1:A:166:ASP:O	1:A:241:GLY:HA2	2.11	0.51
1:B:2:MSE:HE2	1:B:158:VAL:HG11	1.93	0.50
1:C:10:VAL:HG11	1:C:163:THR:O	2.10	0.50
1:C:47:SER:OG	1:C:47:SER:O	2.28	0.50
1:D:50[B]:HIS:HE2	5:D:247:IP8:P1	2.35	0.50
1:B:187:THR:HG23	1:B:210:VAL:HG13	1.94	0.50
1:D:18:ARG:CG	6:D:254:HOH:O	2.59	0.50
1:D:178:ASP:OD1	1:D:178:ASP:C	2.50	0.50
1:A:14:LYS:HE2	1:A:46:GLY:HA2	1.94	0.50
1:B:184:ASP:OD2	1:B:184:ASP:C	2.50	0.50
1:D:168:ILE:HG13	1:D:222:LEU:HD12	1.93	0.49
1:C:10:VAL:HG12	1:C:164:ASP:N	2.28	0.49
1:A:184:ASP:OD1	1:A:245:ARG:CD	2.61	0.49
1:D:176:ASN:OD1	1:D:176:ASN:N	2.45	0.49
1:B:209:MSE:CE	1:B:220:VAL:HG12	2.42	0.49
1:A:62:LYS:HE3	1:A:134:ASP:O	2.13	0.49
1:A:187:THR:C	1:A:188:ASN:OD1	2.51	0.49
1:D:111:TYR:CG	1:D:150:MSE:CE	2.96	0.49
1:D:171:LYS:O	1:D:173:PRO:HD3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:ILE:HD12	1:B:244:ILE:N	2.28	0.49
4:C:246:ATP:O1G	5:C:247:IP8:O2	2.32	0.48
1:B:2:MSE:HG3	1:B:158:VAL:HG12	1.94	0.48
1:D:176:ASN:HB2	1:D:179:ALA:HB2	1.95	0.48
4:C:246:ATP:PG	5:C:247:IP8:P1	3.10	0.48
1:D:171:LYS:HD2	1:D:176:ASN:HD21	1.79	0.47
1:C:29:VAL:HG12	1:C:91:MSE:HG3	1.96	0.47
1:D:15:SER:HB3	6:D:254:HOH:O	2.15	0.47
1:D:171:LYS:HB2	1:D:176:ASN:ND2	2.30	0.47
4:D:246:ATP:PG	5:D:247:IP8:P1	3.10	0.47
1:C:233:ASP:CA	1:C:236:LYS:HD3	2.39	0.47
1:C:166:ASP:HA	1:C:224:ASN:HB2	1.97	0.47
1:C:209:MSE:HE3	1:C:244:ILE:HD11	1.92	0.47
1:C:168:ILE:HD13	1:C:202:ILE:HD13	1.97	0.47
1:C:98:VAL:CB	1:C:114:LEU:CD2	2.92	0.46
1:D:13:ASP:CG	6:D:254:HOH:O	2.54	0.46
1:B:0:THR:HA	1:B:157:ASP:OD2	2.14	0.46
1:B:183:ARG:HH11	1:B:183:ARG:HG3	1.79	0.46
1:D:26:ARG:NH1	1:D:26:ARG:CG	2.75	0.46
1:D:88:MSE:HE2	1:D:95:PRO:HB3	1.97	0.46
1:D:187:THR:HG22	1:D:210:VAL:HG13	1.97	0.46
2:A:246:ADP:O3B	3:A:247:IPE:O2B	2.34	0.46
1:A:49:GLY:C	3:A:247:IPE:H52	2.35	0.46
1:A:171:LYS:HB3	1:A:176:ASN:OD1	2.14	0.46
1:C:13:ASP:HB3	1:C:15:SER:H	1.81	0.46
1:A:185:ILE:HD13	1:A:206:PHE:CE1	2.51	0.45
1:B:207:GLU:O	1:B:211:LYS:HG2	2.15	0.45
1:D:-1:PHE:CD2	1:D:-1:PHE:C	2.89	0.45
1:D:175:ARG:HB2	1:D:176:ASN:OD1	2.17	0.45
1:B:227:HIS:O	1:B:230:ARG:HG3	2.17	0.45
1:D:18:ARG:HD2	1:D:20:ALA:O	2.17	0.44
1:B:166:ASP:HA	1:B:224:ASN:HB2	1.99	0.44
1:C:35:ILE:HD11	1:C:231:ILE:HG21	1.99	0.44
1:B:114:LEU:C	1:B:114:LEU:HD23	2.37	0.44
1:B:112:THR:N	1:B:113:PRO:CD	2.80	0.44
1:D:31:VAL:HG21	1:D:228:PRO:HB2	1.99	0.44
1:D:174:LYS:HB3	1:D:174:LYS:HE2	1.79	0.44
1:C:98:VAL:HG11	1:C:114:LEU:HD23	1.99	0.44
1:A:14:LYS:CD	6:A:287:HOH:O	2.64	0.44
1:C:63:ASN:HB2	1:C:64:PRO:CD	2.48	0.44
1:A:202:ILE:CA	6:A:289:HOH:O	2.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:THR:HG22	1:C:244:ILE:CG2	2.47	0.44
1:D:150:MSE:CE	1:D:153:LEU:HD12	2.48	0.44
1:C:161:PHE:HB2	1:C:222:LEU:HD12	1.99	0.43
1:D:96:ILE:O	1:D:96:ILE:HG23	2.19	0.43
2:A:246:ADP:O3B	3:A:247:IPE:O1B	2.35	0.43
1:B:183:ARG:HG3	1:B:183:ARG:NH1	2.34	0.43
1:A:88:MSE:HE2	1:A:95:PRO:HB3	2.00	0.43
1:B:10:VAL:HG22	1:B:164:ASP:CA	2.48	0.43
1:B:221:TYR:CE2	1:B:243:VAL:HG13	2.54	0.43
1:A:68:ILE:HD13	1:C:94:ARG:HG2	2.01	0.43
1:A:185:ILE:HD13	1:A:206:PHE:HE1	1.84	0.43
1:D:150:MSE:HA	1:D:150:MSE:HE2	1.99	0.43
1:D:151:ALA:CB	1:D:212:MSE:HE3	2.49	0.43
1:D:168:ILE:HD12	1:D:242:THR:HG21	2.01	0.42
1:C:5:LYS:HA	1:C:42:VAL:O	2.19	0.42
1:D:150:MSE:HE1	1:D:153:LEU:CD1	2.49	0.42
1:C:111:TYR:CD1	1:C:150:MSE:CE	3.02	0.42
1:A:166:ASP:HA	1:A:224:ASN:HB2	2.02	0.42
1:D:111:TYR:CD2	1:D:150:MSE:CE	3.03	0.42
1:D:23:TYR:CD2	1:D:23:TYR:C	2.93	0.42
1:A:176:ASN:N	1:A:176:ASN:ND2	2.68	0.42
1:D:221:TYR:CE2	1:D:243:VAL:CG2	3.00	0.42
1:C:64:PRO:O	1:C:68:ILE:HG12	2.20	0.41
1:B:0:THR:HB	1:B:157:ASP:CB	2.50	0.41
1:A:176:ASN:ND2	1:A:176:ASN:H	2.18	0.41
1:B:10:VAL:HG22	1:B:164:ASP:N	2.35	0.41
1:D:158:VAL:HG22	1:D:218:ASN:HB2	2.02	0.41
1:C:187:THR:HG23	1:C:245:ARG:O	2.21	0.41
1:C:91:MSE:N	1:C:91:MSE:HE2	2.35	0.41
1:A:82:LEU:CD1	1:C:82:LEU:HD12	2.50	0.41
1:A:17:TYR:CE1	1:A:52:LYS:HG2	2.56	0.41
1:A:86:ASP:OD1	1:C:75:ARG:HD3	2.20	0.41
2:B:246:ADP:N3	2:B:246:ADP:H2'	2.35	0.41
1:A:206:PHE:HA	1:A:209:MSE:HE3	2.02	0.41
1:B:106:ASP:O	1:B:106:ASP:OD1	2.38	0.41
1:D:87:ALA:O	1:D:91:MSE:HG2	2.20	0.41
1:B:77:MSE:HE2	1:B:128:GLY:O	2.20	0.41
1:B:75:ARG:HD2	1:B:75:ARG:C	2.41	0.41
1:D:98:VAL:HG11	1:D:114:LEU:CD1	2.50	0.41
1:B:207:GLU:O	1:B:211:LYS:CG	2.70	0.40
1:A:68:ILE:HA	1:A:68:ILE:HD13	1.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LEU:HD12	1:C:82:LEU:CD1	2.52	0.40
1:C:150:MSE:HA	1:C:150:MSE:HE2	2.03	0.40
1:C:213:LYS:HA	1:C:213:LYS:HD2	1.86	0.40
1:C:151:ALA:HB2	1:C:212:MSE:HE3	2.03	0.40
1:A:202:ILE:HG23	1:A:203:GLY:N	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	229/249 (92%)	223 (97%)	6 (3%)	0	100	100
1	B	230/249 (92%)	227 (99%)	3 (1%)	0	100	100
1	C	234/249 (94%)	229 (98%)	5 (2%)	0	100	100
1	D	231/249 (93%)	214 (93%)	15 (6%)	2 (1%)	21	12
All	All	924/996 (93%)	893 (97%)	29 (3%)	2 (0%)	52	47

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	236	LYS
1	D	172	ASP

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	199/201 (99%)	194 (98%)	5 (2%)	55	53
1	B	199/201 (99%)	196 (98%)	3 (2%)	72	73
1	C	203/201 (101%)	196 (97%)	7 (3%)	44	39
1	D	201/201 (100%)	193 (96%)	8 (4%)	38	31
All	All	802/804 (100%)	779 (97%)	23 (3%)	50	46

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ILE
1	A	135	GLU
1	A	176	ASN
1	A	178	ASP
1	A	184	ASP
1	B	100	ILE
1	B	178	ASP
1	B	205	LYS
1	C	18	ARG
1	C	54	MSE
1	C	79	ASN
1	C	114	LEU
1	C	137	SER
1	C	176	ASN
1	C	187	THR
1	D	10	VAL
1	D	26	ARG
1	D	27	SER
1	D	176	ASN
1	D	178	ASP
1	D	204	LYS
1	D	214	SER
1	D	215	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	246	-	22,29,29	1.13	3 (13%)	27,45,45	2.13	6 (22%)
3	IPE	A	247	-	10,13,13	0.88	0	14,19,19	2.09	4 (28%)
2	ADP	B	246	-	22,29,29	1.19	2 (9%)	27,45,45	1.95	6 (22%)
3	IPE	B	247	-	10,13,13	0.51	0	14,19,19	2.01	6 (42%)
4	ATP	C	246	-	24,33,33	1.08	1 (4%)	31,52,52	1.58	3 (9%)
5	IP8	C	247	-	8,9,9	2.76	4 (50%)	8,12,12	2.88	6 (75%)
4	ATP	D	246	-	24,33,33	1.20	2 (8%)	31,52,52	1.71	6 (19%)
5	IP8	D	247	-	8,9,9	2.39	3 (37%)	8,12,12	2.05	4 (50%)

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	246	-	-	0/12/32/32	0/3/3/3
3	IPE	A	247	-	-	0/13/13/13	0/0/0/0
2	ADP	B	246	-	-	0/12/32/32	0/3/3/3
3	IPE	B	247	-	-	0/13/13/13	0/0/0/0
4	ATP	C	246	-	-	0/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	IP8	C	247	-	-	0/7/7/7	0/0/0/0
4	ATP	D	246	-	-	0/18/38/38	0/3/3/3
5	IP8	D	247	-	-	0/7/7/7	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	247	IP8	C4-C3	-3.74	1.31	1.48
5	D	247	IP8	C4-C3	-3.50	1.32	1.48
2	A	246	ADP	C2-N3	2.03	1.35	1.32
5	D	247	IP8	P1-O1	2.24	1.67	1.60
5	C	247	IP8	C2-C1	2.28	1.57	1.51
2	A	246	ADP	O4'-C1'	2.28	1.44	1.41
2	B	246	ADP	C5-C4	2.69	1.46	1.40
4	D	246	ATP	O4'-C1'	2.85	1.44	1.41
5	C	247	IP8	P1-O1	2.96	1.70	1.60
2	A	246	ADP	C5-C4	3.02	1.47	1.40
2	B	246	ADP	O4'-C1'	3.20	1.45	1.41
4	D	246	ATP	C5-C4	3.43	1.48	1.40
4	C	246	ATP	C5-C4	3.64	1.48	1.40
5	D	247	IP8	C5-C3	4.88	1.48	1.33
5	C	247	IP8	C5-C3	5.44	1.50	1.33

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	246	ADP	N3-C2-N1	-7.77	122.94	128.89
4	D	246	ATP	N3-C2-N1	-6.13	124.20	128.89
2	B	246	ADP	N3-C2-N1	-5.97	124.32	128.89
5	C	247	IP8	O3-P1-O1	-4.96	92.27	106.56
4	C	246	ATP	N3-C2-N1	-4.80	125.22	128.89
2	B	246	ADP	C4'-O4'-C1'	-4.32	104.97	109.72
4	C	246	ATP	O3A-PA-O5'	-4.26	91.64	102.94
2	A	246	ADP	C4'-O4'-C1'	-4.06	105.26	109.72
3	A	247	IPE	PA-O3A-PB	-3.99	119.28	132.67
3	B	247	IPE	O3A-PA-O1	-3.39	93.94	102.94
4	D	246	ATP	O3A-PA-O5'	-3.01	94.95	102.94
5	C	247	IP8	C4-C3-C5	-2.99	114.18	121.72
4	C	246	ATP	C4-C5-N7	-2.89	106.82	109.48
2	B	246	ADP	C2'-C1'-N9	-2.84	109.96	114.29
3	A	247	IPE	O3B-PB-O3A	-2.72	92.77	105.09
2	A	246	ADP	PA-O3A-PB	-2.72	123.56	132.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	247	IP8	O3-P1-O1	-2.62	99.01	106.56
2	B	246	ADP	PA-O3A-PB	-2.59	123.97	132.67
4	D	246	ATP	C4-C5-N7	-2.59	107.10	109.48
3	B	247	IPE	O1B-PB-O3A	-2.53	93.61	105.09
5	D	247	IP8	C4-C3-C5	-2.41	115.65	121.72
4	D	246	ATP	PA-O3A-PB	-2.30	126.27	132.73
3	A	247	IPE	O1B-PB-O3A	-2.22	95.02	105.09
2	A	246	ADP	O5'-C5'-C4'	-2.20	100.99	109.12
2	B	246	ADP	C4-C5-N7	-2.18	107.48	109.48
3	B	247	IPE	PA-O3A-PB	-2.15	125.47	132.67
2	A	246	ADP	C4-C5-N7	-2.07	107.57	109.48
3	B	247	IPE	O2A-PA-O3A	2.11	114.66	105.09
3	B	247	IPE	O3B-PB-O1B	2.12	115.45	107.38
4	D	246	ATP	O2B-PB-O1B	2.28	124.91	112.53
4	D	246	ATP	C2'-C1'-N9	2.32	117.84	114.29
5	C	247	IP8	O1-P1-O	2.35	113.12	107.14
5	D	247	IP8	O1-P1-O	2.50	113.51	107.14
2	B	246	ADP	O2A-PA-O3A	2.62	116.97	105.09
5	C	247	IP8	O2-P1-O3	2.71	117.71	107.38
2	A	246	ADP	O3B-PB-O2B	2.74	117.83	107.38
5	D	247	IP8	O2-P1-O3	2.78	117.97	107.38
5	C	247	IP8	O1-C1-C2	2.93	122.69	108.63
5	C	247	IP8	O3-P1-O	3.05	120.41	110.58
3	B	247	IPE	O1B-PB-O2B	4.24	124.22	110.58
3	A	247	IPE	O3B-PB-O1B	4.30	123.76	107.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	246	ADP	4	0
3	A	247	IPE	5	0
2	B	246	ADP	3	0
3	B	247	IPE	2	0
4	C	246	ATP	6	0
5	C	247	IP8	13	0
4	D	246	ATP	5	0
5	D	247	IP8	11	0

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 ⓘ

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	221/249 (88%)	-0.12	6 (2%) 58 61	15, 25, 50, 73	2 (0%)
1	B	222/249 (89%)	0.11	8 (3%) 46 50	16, 26, 50, 75	1 (0%)
1	C	225/249 (90%)	0.02	10 (4%) 38 42	15, 24, 51, 60	2 (0%)
1	D	222/249 (89%)	0.63	29 (13%) 5 6	18, 32, 69, 84	1 (0%)
All	All	890/996 (89%)	0.16	53 (5%) 25 29	15, 26, 58, 84	6 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	185	ILE	8.1
1	C	48	PHE	7.0
1	D	177	PRO	5.8
1	B	188	ASN	5.7
1	D	175	ARG	5.2
1	A	177	PRO	5.2
1	B	187	THR	4.8
1	A	178	ASP	4.4
1	A	202	ILE	4.2
1	D	178	ASP	4.0
1	B	0	THR	3.9
1	D	187	THR	3.9
1	A	175	ARG	3.7
1	D	48	PHE	3.6
1	B	201	GLY	3.5
1	D	206	PHE	3.4
1	D	188	ASN	3.3
1	D	236	LYS	3.2
1	D	186	ASP	3.1
1	D	239	PHE	3.1
1	C	187	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	175	ARG	3.0
1	B	202	ILE	3.0
1	C	177	PRO	2.9
1	D	237	GLU	2.9
1	C	47	SER	2.8
1	D	217	LYS	2.8
1	D	-1	PHE	2.8
1	D	179	ALA	2.7
1	D	181	LEU	2.7
1	B	186	ASP	2.7
1	D	184	ASP	2.6
1	A	206	PHE	2.6
1	D	240	ILE	2.6
1	D	243	VAL	2.5
1	D	218	ASN	2.5
1	C	178	ASP	2.5
1	D	171	LYS	2.5
1	D	170	SER	2.5
1	D	235	GLY	2.5
1	C	136	HIS	2.5
1	D	245	ARG	2.4
1	D	102	ALA	2.3
1	D	34	GLY	2.3
1	D	183	ARG	2.3
1	C	14	LYS	2.2
1	D	202	ILE	2.2
1	D	238	SER	2.2
1	C	19	THR	2.2
1	C	238	SER	2.2
1	B	135	GLU	2.1
1	B	102	ALA	2.1
1	A	176	ASN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	IP8	C	247	10/10	0.84	0.18	3.00	30,36,41,52	1
5	IP8	D	247	10/10	0.91	0.14	1.50	30,36,42,52	2
3	IPE	A	247	14/14	0.94	0.11	0.62	29,34,41,41	0
4	ATP	C	246	31/31	0.97	0.10	-0.33	25,40,50,51	0
3	IPE	B	247	14/14	0.97	0.09	-0.47	24,31,39,40	0
4	ATP	D	246	31/31	0.97	0.10	-0.63	27,40,52,55	0
2	ADP	A	246	27/27	0.97	0.08	-0.65	27,37,43,47	1
2	ADP	B	246	27/27	0.98	0.07	-0.86	23,27,35,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.