



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:16 PM GMT

PDB ID : 1O93
Title : METHIONINE ADENOSYLTRANSFERASE COMPLEXED WITH ATP
AND A L-METHIONINE ANALOGOUS
Authors : Gonzalez, B.; Pajares, M.A.; Hermoso, J.A.; Sanz-Aparicio, J.
Deposited on : 2002-12-10
Resolution : 3.49 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

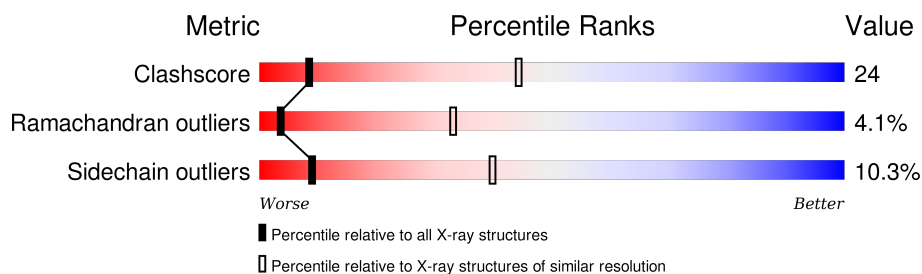
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.49 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))
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	396	
1	B	396	

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
3	PO4	B	1399	-	-	X	-
6	ATP	B	1397	-	-	X	-

2 Entry composition [i](#)

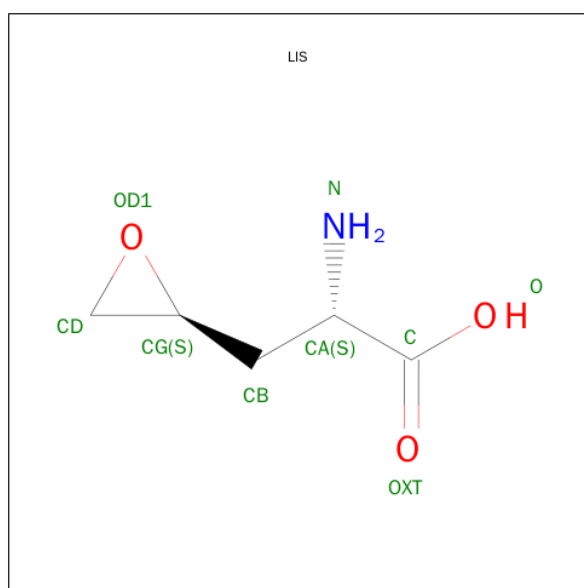
There are 7 unique types of molecules in this entry. The entry contains 5932 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 S-ADENOSYLMETHIONINE SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2846	1803	491	536	16			
1	B	368	Total	C	N	O	S	0	0	0
			2846	1803	491	536	16			

- Molecule 2 is (2S,4S)-2-AMINO-4,5-EPOXIPENTANOIC ACID (three-letter code: LIS) (formula: C₅H₉NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	5	1	3		

- Molecule 3 is PHOSPHATE ION (three-letter code: 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		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	K	0	0
			1	1		
5	A	1	Total	K	0	0
			1	1		

- Molecule 6 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 7 is water.

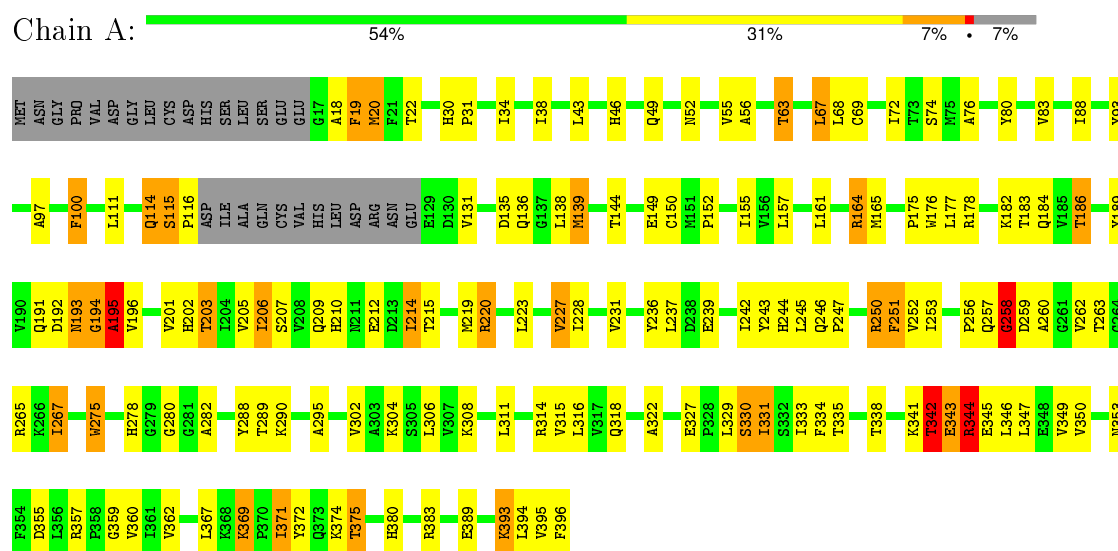
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	102	Total O 102 102	0	0
7	B	83	Total O 83 83	0	0

3 Residue-property plots

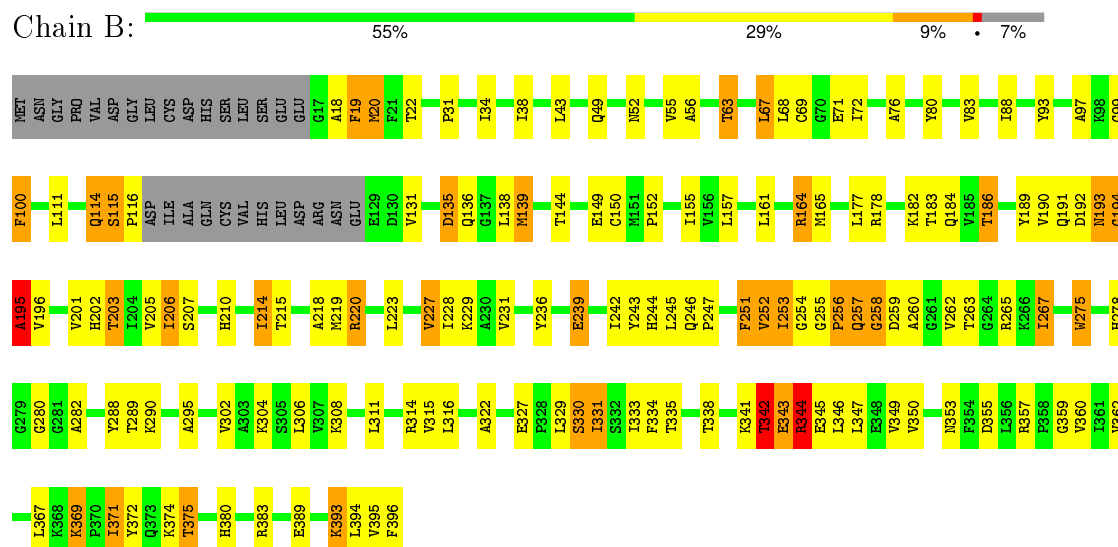
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

• Molecule 1: S-ADENOSYLMETHIONINE SYNTHETASE



• Molecule 1: S-ADENOSYLMETHIONINE SYNTHETASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	115.20 Å 115.20 Å 160.89 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.49	Depositor
% Data completeness (in resolution range)	97.3 (8.00-3.49)	Depositor
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.224 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5932	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, MG, K, ATP, LIS

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	0.47	1/2901 (0.0%)	0.83	9/3926 (0.2%)
1	B	0.39	0/2901	0.72	5/3926 (0.1%)
All	All	0.43	1/5802 (0.0%)	0.78	14/7852 (0.2%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	258	GLY	C-N	12.57	1.62	1.34

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	258	GLY	O-C-N	-20.19	90.40	122.70
1	A	258	GLY	CA-C-N	-8.44	98.64	117.20
1	A	195	ALA	O-C-N	6.78	133.56	122.70
1	A	344	ARG	NE-CZ-NH2	6.58	123.59	120.30
1	B	178	ARG	NE-CZ-NH2	6.52	123.56	120.30
1	A	178	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	A	250	ARG	NE-CZ-NH2	6.40	123.50	120.30
1	B	195	ALA	O-C-N	6.32	132.81	122.70
1	B	20	MET	CG-SD-CE	6.20	110.12	100.20
1	A	20	MET	CG-SD-CE	6.12	110.00	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	220	ARG	NE-CZ-NH2	5.89	123.24	120.30
1	A	220	ARG	NE-CZ-NH2	5.78	123.19	120.30
1	A	195	ALA	CA-C-N	-5.29	105.56	117.20
1	B	195	ALA	CA-C-N	-5.17	105.84	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	258	GLY	Mainchain

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	2846	0	2852	137	0
1	B	2846	0	2852	147	0
2	A	9	0	7	1	0
3	A	5	0	0	1	0
3	B	5	0	0	2	0
4	A	3	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	31	0	12	13	0
7	A	102	0	0	3	0
7	B	83	0	0	4	0
All	All	5932	0	5723	275	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 24.

All (275) 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:A:19:PHE:CE1	1:A:189:TYR:HB2	1.80	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:PHE:CE1	1:B:189:TYR:HB2	1.85	1.10
1:B:290:LYS:NZ	6:B:1397:ATP:C2	2.23	1.06
1:B:290:LYS:NZ	6:B:1397:ATP:N3	2.06	1.02
1:A:19:PHE:HE1	1:A:189:TYR:HB2	1.12	1.01
1:A:267:ILE:HD11	1:A:282:ALA:HA	1.42	1.01
1:B:267:ILE:HD11	1:B:282:ALA:HA	1.41	1.01
1:A:19:PHE:HE1	1:A:189:TYR:CB	1.77	0.96
1:B:100:PHE:HD2	1:B:256:PRO:HG2	1.31	0.96
1:B:100:PHE:CD2	1:B:256:PRO:HG2	2.02	0.95
1:A:100:PHE:HD2	1:A:256:PRO:HG2	1.34	0.93
1:B:19:PHE:HE1	1:B:189:TYR:HB2	1.20	0.92
1:A:262:VAL:HG11	1:B:262:VAL:HG11	1.53	0.91
1:B:19:PHE:HD1	1:B:19:PHE:O	1.53	0.90
1:A:100:PHE:CD2	1:A:256:PRO:HG2	2.07	0.90
1:B:19:PHE:HE1	1:B:189:TYR:CB	1.87	0.87
1:A:203:THR:HB	1:A:242:ILE:HB	1.59	0.85
1:B:203:THR:HB	1:B:242:ILE:HB	1.59	0.84
1:B:278:HIS:HD2	1:B:280:GLY:H	1.27	0.83
1:B:205:VAL:HG22	1:B:244:HIS:HB2	1.63	0.80
1:A:278:HIS:HD2	1:A:280:GLY:H	1.26	0.79
1:A:205:VAL:HG22	1:A:244:HIS:HB2	1.64	0.79
1:A:19:PHE:HD1	1:A:19:PHE:O	1.66	0.78
1:B:290:LYS:CE	6:B:1397:ATP:N3	2.47	0.77
1:B:210:HIS:CE1	1:B:251:PHE:H	2.03	0.77
1:A:257:GLN:HG3	1:A:258:GLY:N	2.02	0.74
1:B:19:PHE:CD1	1:B:19:PHE:C	2.59	0.74
1:A:258:GLY:CA	6:B:1397:ATP:O2A	2.36	0.74
1:B:55:VAL:HA	1:B:72:ILE:HG22	1.69	0.73
1:A:55:VAL:HA	1:A:72:ILE:HG22	1.70	0.72
1:A:257:GLN:CG	1:A:258:GLY:N	2.52	0.71
1:B:19:PHE:CD1	1:B:19:PHE:O	2.42	0.71
1:A:19:PHE:CD1	1:A:189:TYR:HB2	2.27	0.70
1:A:322:ALA:HB2	1:A:329:LEU:HD22	1.73	0.70
1:A:267:ILE:HG23	1:A:278:HIS:CE1	2.26	0.70
1:A:31:PRO:HG3	1:A:256:PRO:HA	1.74	0.69
1:A:19:PHE:CE1	1:A:189:TYR:CB	2.62	0.69
1:B:19:PHE:HD1	1:B:19:PHE:C	1.95	0.69
1:B:267:ILE:HG23	1:B:278:HIS:CE1	2.26	0.69
1:B:322:ALA:HB2	1:B:329:LEU:HD22	1.74	0.69
1:A:182:LYS:HE3	7:A:2001:HOH:O	1.93	0.68
1:A:19:PHE:CD1	1:A:19:PHE:C	2.66	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:GLY:O	1:A:195:ALA:HB2	1.95	0.67
1:A:186:THR:CG2	1:A:203:THR:HG23	2.25	0.67
1:B:186:THR:CG2	1:B:203:THR:HG23	2.25	0.67
1:B:72:ILE:HG12	1:B:111:LEU:HD11	1.76	0.67
1:B:194:GLY:O	1:B:195:ALA:HB2	1.95	0.67
1:A:342:THR:O	1:A:345:GLU:HB2	1.94	0.66
1:A:72:ILE:HG12	1:A:111:LEU:HD11	1.78	0.66
1:B:99:GLY:HA3	1:B:255:GLY:HA3	1.77	0.66
1:A:329:LEU:O	1:A:330:SER:HB3	1.96	0.65
1:A:138:LEU:HD22	1:B:22:THR:H	1.62	0.65
1:B:19:PHE:CD1	1:B:189:TYR:HB2	2.32	0.65
1:A:257:GLN:CG	1:A:258:GLY:H	2.09	0.65
1:B:220:ARG:HD2	7:B:2045:HOH:O	1.97	0.65
1:B:329:LEU:O	1:B:330:SER:HB3	1.97	0.64
1:A:184:GLN:OE1	1:B:136:GLN:HB2	1.97	0.64
1:B:290:LYS:HZ2	6:B:1397:ATP:H2	1.45	0.64
1:A:335:THR:HG22	1:A:338:THR:H	1.62	0.63
1:A:210:HIS:CE1	1:A:251:PHE:H	2.17	0.63
1:B:335:THR:HG22	1:B:338:THR:H	1.62	0.63
1:A:353:ASN:HD22	1:A:394:LEU:HD22	1.64	0.63
1:B:56:ALA:CB	6:B:1397:ATP:C5	2.82	0.63
1:B:19:PHE:CE1	1:B:189:TYR:CB	2.68	0.63
1:B:353:ASN:HD22	1:B:394:LEU:HD22	1.64	0.63
1:A:19:PHE:CD1	1:A:19:PHE:O	2.50	0.62
1:B:19:PHE:HE1	1:B:189:TYR:CG	2.17	0.62
1:A:333:ILE:O	1:A:343:GLU:HG2	2.00	0.62
1:B:186:THR:HG23	1:B:203:THR:HG23	1.82	0.61
1:B:333:ILE:O	1:B:343:GLU:HG2	2.00	0.61
1:A:19:PHE:HE1	1:A:189:TYR:CG	2.17	0.61
1:A:186:THR:HG23	1:A:203:THR:HG23	1.83	0.61
1:A:334:PHE:HZ	1:B:19:PHE:HD2	1.48	0.60
1:B:56:ALA:HB2	6:B:1397:ATP:C4	2.38	0.59
1:A:258:GLY:C	6:B:1397:ATP:O2A	2.41	0.59
1:B:31:PRO:HG3	1:B:256:PRO:HA	1.84	0.59
1:A:250:ARG:O	2:A:1398:LIS:HDC2	2.02	0.59
1:A:345:GLU:O	1:A:349:VAL:HG23	2.03	0.59
1:B:56:ALA:H	1:B:72:ILE:HG22	1.68	0.58
1:B:257:GLN:O	1:B:260:ALA:N	2.35	0.58
1:A:56:ALA:H	1:A:72:ILE:HG22	1.68	0.58
1:B:344:ARG:NE	1:B:347:LEU:HD23	2.18	0.58
1:B:228:ILE:HG21	1:B:243:TYR:HE2	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:LYS:HE2	1:B:184:GLN:HE21	1.69	0.58
1:A:344:ARG:NE	1:A:347:LEU:HD23	2.18	0.57
1:B:219:MET:HE1	1:B:246:GLN:HE21	1.69	0.57
1:A:258:GLY:HA2	6:B:1397:ATP:O2A	2.03	0.57
1:A:228:ILE:HG21	1:A:243:TYR:HE2	1.70	0.57
1:B:223:LEU:O	1:B:227:VAL:HG22	2.06	0.56
1:A:67:LEU:HD22	1:A:69:CYS:SG	2.45	0.56
1:B:369:LYS:O	1:B:371:ILE:HD13	2.05	0.56
1:B:67:LEU:HD22	1:B:69:CYS:SG	2.46	0.56
1:B:100:PHE:HD2	1:B:256:PRO:CG	2.14	0.55
1:B:114:GLN:HE21	1:B:114:GLN:N	2.05	0.55
1:A:192:ASP:O	1:A:194:GLY:N	2.39	0.55
1:B:282:ALA:HB2	3:B:1399:PO4:O3	2.05	0.55
1:B:393:LYS:HD2	1:B:393:LYS:H	1.70	0.55
1:A:393:LYS:HD2	1:A:393:LYS:H	1.71	0.55
1:A:369:LYS:O	1:A:371:ILE:HD13	2.06	0.55
1:A:375:THR:HG23	1:A:380:HIS:NE2	2.22	0.55
1:A:223:LEU:O	1:A:227:VAL:HG22	2.07	0.55
1:B:375:THR:HG23	1:B:380:HIS:NE2	2.22	0.55
1:B:72:ILE:HD11	1:B:80:TYR:OH	2.07	0.55
1:B:56:ALA:CB	6:B:1397:ATP:C4	2.90	0.55
1:B:246:GLN:N	1:B:247:PRO:HD3	2.22	0.54
1:A:72:ILE:HD11	1:A:80:TYR:OH	2.07	0.54
1:B:383:ARG:O	1:B:389:GLU:HG3	2.08	0.54
1:B:71:GLU:HB3	6:B:1397:ATP:HN61	1.73	0.54
1:A:114:GLN:N	1:A:114:GLN:HE21	2.05	0.54
1:B:257:GLN:O	1:B:259:ASP:N	2.41	0.54
1:A:246:GLN:N	1:A:247:PRO:HD3	2.23	0.54
1:B:345:GLU:O	1:B:349:VAL:HG23	2.07	0.54
1:A:186:THR:HG21	1:A:203:THR:HG23	1.91	0.53
1:A:136:GLN:NE2	1:A:322:ALA:HA	2.22	0.53
1:A:161:LEU:HD23	1:A:183:THR:HG23	1.91	0.53
1:A:219:MET:CE	1:A:246:GLN:HE21	2.22	0.53
1:A:93:TYR:CZ	1:A:256:PRO:HG3	2.44	0.53
1:A:383:ARG:O	1:A:389:GLU:HG3	2.09	0.52
1:A:334:PHE:CZ	1:B:19:PHE:HD2	2.27	0.52
1:A:257:GLN:HG2	1:A:258:GLY:H	1.73	0.52
1:A:207:SER:HA	1:A:246:GLN:OE1	2.09	0.52
1:B:186:THR:HG21	1:B:203:THR:HG23	1.92	0.52
1:B:219:MET:CE	1:B:246:GLN:HE21	2.22	0.52
1:B:72:ILE:CG1	1:B:111:LEU:HD11	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:LEU:HD23	1:B:183:THR:HG23	1.91	0.52
1:B:52:ASN:HA	1:B:288:TYR:OH	2.10	0.52
1:B:165:MET:SD	1:B:206:ILE:HD11	2.50	0.52
1:A:52:ASN:HA	1:A:288:TYR:OH	2.10	0.52
1:A:316:LEU:HB3	1:A:334:PHE:HB2	1.92	0.51
1:A:202:HIS:ND1	1:A:203:THR:HG22	2.26	0.51
1:B:275:TRP:HE1	1:B:314:ARG:HH21	1.58	0.51
1:B:316:LEU:HB3	1:B:334:PHE:HB2	1.92	0.51
1:A:219:MET:HE1	1:A:246:GLN:HE21	1.75	0.51
1:A:201:VAL:HG23	1:A:236:TYR:HB3	1.93	0.51
1:B:207:SER:HA	1:B:246:GLN:OE1	2.11	0.51
1:B:202:HIS:ND1	1:B:203:THR:HG22	2.26	0.50
1:B:210:HIS:HE1	1:B:251:PHE:H	1.54	0.50
1:B:192:ASP:O	1:B:194:GLY:N	2.44	0.50
1:A:165:MET:SD	1:A:206:ILE:HD11	2.50	0.50
1:A:275:TRP:HE1	1:A:314:ARG:HH21	1.58	0.50
1:A:258:GLY:O	1:A:259:ASP:OD1	2.29	0.50
1:A:22:THR:H	1:B:138:LEU:HD22	1.77	0.50
1:B:227:VAL:O	1:B:231:VAL:HG22	2.12	0.50
1:A:257:GLN:O	1:A:259:ASP:N	2.44	0.50
1:B:201:VAL:HG23	1:B:236:TYR:HB3	1.93	0.50
1:B:282:ALA:CB	3:B:1399:PO4:O3	2.60	0.49
1:A:22:THR:OG1	1:A:186:THR:HB	2.12	0.49
1:A:346:LEU:O	1:A:350:VAL:HG23	2.12	0.49
1:B:220:ARG:HG2	1:B:245:LEU:HB2	1.93	0.49
1:A:72:ILE:CG1	1:A:111:LEU:HD11	2.42	0.49
1:A:227:VAL:O	1:A:231:VAL:HG22	2.12	0.49
1:A:97:ALA:O	1:A:253:ILE:HG23	2.12	0.49
1:B:346:LEU:O	1:B:350:VAL:HG23	2.13	0.49
1:B:302:VAL:O	1:B:306:LEU:HD13	2.12	0.49
1:B:290:LYS:HE2	6:B:1397:ATP:H1'	1.95	0.49
1:B:43:LEU:HD13	1:B:55:VAL:HB	1.95	0.49
1:B:22:THR:OG1	1:B:186:THR:HB	2.12	0.49
1:B:136:GLN:NE2	1:B:322:ALA:HA	2.27	0.48
1:A:353:ASN:ND2	1:A:394:LEU:HA	2.28	0.48
1:A:182:LYS:NZ	3:A:1400:PO4:O3	2.34	0.48
1:A:175:PRO:HA	7:A:2046:HOH:O	2.12	0.48
1:B:252:VAL:HG12	1:B:253:ILE:HD13	1.95	0.48
1:B:97:ALA:O	1:B:253:ILE:HG23	2.13	0.48
1:A:52:ASN:HA	1:A:288:TYR:HH	1.78	0.48
1:B:152:PRO:HB2	1:B:155:ILE:HG22	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:PHE:HD2	1:B:334:PHE:HZ	1.60	0.48
1:A:395:VAL:O	1:A:395:VAL:HG12	2.14	0.48
1:A:152:PRO:HB2	1:A:155:ILE:HG22	1.96	0.48
1:A:362:VAL:HA	1:A:367:LEU:HD12	1.96	0.48
1:B:362:VAL:HA	1:B:367:LEU:HD12	1.95	0.48
1:B:34:ILE:O	1:B:38:ILE:HG13	2.14	0.48
1:A:220:ARG:HG2	1:A:245:LEU:HB2	1.96	0.47
1:A:144:THR:O	1:A:150:CYS:HA	2.14	0.47
1:A:43:LEU:HD13	1:A:55:VAL:HB	1.95	0.47
1:B:353:ASN:ND2	1:B:394:LEU:HA	2.30	0.47
1:B:395:VAL:O	1:B:395:VAL:HG12	2.14	0.47
1:B:161:LEU:O	1:B:165:MET:HG3	2.14	0.47
1:A:302:VAL:O	1:A:306:LEU:HD13	2.13	0.47
1:A:34:ILE:O	1:A:38:ILE:HG13	2.13	0.47
1:B:359:GLY:O	1:B:362:VAL:HG22	2.14	0.47
1:B:214:ILE:HD12	1:B:219:MET:CG	2.45	0.47
1:B:372:TYR:O	1:B:375:THR:HG22	2.14	0.47
1:A:214:ILE:HD12	1:A:219:MET:CG	2.45	0.46
1:B:219:MET:HE1	1:B:246:GLN:NE2	2.31	0.46
1:A:372:TYR:O	1:A:375:THR:HG22	2.15	0.46
1:B:144:THR:O	1:B:150:CYS:HA	2.15	0.46
1:B:344:ARG:HD2	1:B:344:ARG:HA	1.64	0.46
1:B:80:TYR:HA	1:B:83:VAL:HG22	1.96	0.46
1:B:165:MET:HE2	1:B:177:LEU:HD11	1.98	0.46
1:A:311:LEU:HD22	1:A:341:LYS:HD2	1.97	0.46
1:A:30:HIS:HE2	6:B:1397:ATP:PG	2.38	0.46
1:A:161:LEU:O	1:A:165:MET:HG3	2.15	0.46
1:A:80:TYR:HA	1:A:83:VAL:HG22	1.96	0.45
1:B:257:GLN:C	1:B:259:ASP:N	2.69	0.45
1:A:192:ASP:O	1:A:193:ASN:C	2.54	0.45
1:B:259:ASP:O	1:B:260:ALA:HB3	2.16	0.45
1:B:311:LEU:HD22	1:B:341:LYS:HD2	1.97	0.45
1:A:371:ILE:HB	1:A:374:LYS:HE2	1.99	0.45
1:B:165:MET:CE	1:B:206:ILE:HD11	2.47	0.45
1:B:194:GLY:O	1:B:195:ALA:CB	2.62	0.45
1:A:165:MET:CE	1:A:206:ILE:HD11	2.46	0.45
1:B:72:ILE:CD1	1:B:111:LEU:HD11	2.47	0.45
1:A:395:VAL:O	1:A:396:PHE:HB2	2.17	0.45
1:B:327:GLU:HA	1:B:357:ARG:HG2	1.99	0.45
1:A:329:LEU:O	1:B:20:MET:SD	2.75	0.45
1:A:258:GLY:C	1:A:259:ASP:OD1	2.55	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:331:ILE:HG23	1:B:347:LEU:HD11	1.99	0.45
1:A:165:MET:HE2	1:A:177:LEU:HD11	1.99	0.45
1:B:371:ILE:O	1:B:374:LYS:HG2	2.16	0.44
1:A:331:ILE:HG23	1:A:347:LEU:HD11	1.98	0.44
1:A:344:ARG:HA	1:A:344:ARG:HD2	1.62	0.44
1:A:327:GLU:HA	1:A:357:ARG:HG2	1.99	0.44
1:A:371:ILE:O	1:A:374:LYS:HG2	2.17	0.44
1:A:49:GLN:HB2	1:A:76:ALA:HB1	2.00	0.44
1:B:395:VAL:O	1:B:396:PHE:HB2	2.18	0.44
1:B:149:GLU:O	1:B:150:CYS:HB2	2.18	0.44
1:A:72:ILE:CD1	1:A:111:LEU:HD11	2.48	0.44
1:A:359:GLY:O	1:A:362:VAL:HG22	2.17	0.44
1:A:20:MET:HG2	7:B:2068:HOH:O	2.18	0.44
1:B:262:VAL:CG1	1:B:263:THR:N	2.81	0.44
1:B:239:GLU:HG2	7:B:2051:HOH:O	2.18	0.44
1:B:371:ILE:HB	1:B:374:LYS:HE2	1.99	0.44
1:A:149:GLU:O	1:A:150:CYS:HB2	2.18	0.43
1:B:229:LYS:HE3	7:B:2048:HOH:O	2.18	0.43
1:B:267:ILE:HD13	1:B:267:ILE:HA	1.77	0.43
1:A:194:GLY:O	1:A:195:ALA:CB	2.62	0.43
1:B:342:THR:O	1:B:345:GLU:HB2	2.18	0.43
1:A:290:LYS:HE3	7:A:2075:HOH:O	2.17	0.43
1:B:49:GLN:HB2	1:B:76:ALA:HB1	2.00	0.43
1:B:192:ASP:O	1:B:193:ASN:C	2.57	0.43
1:A:115:SER:H	1:A:116:PRO:HD2	1.84	0.43
1:B:115:SER:H	1:B:116:PRO:HD2	1.84	0.43
1:A:355:ASP:OD2	1:A:360:VAL:HG21	2.19	0.42
1:A:237:LEU:HD23	1:A:237:LEU:HA	1.83	0.42
1:B:227:VAL:CG2	1:B:228:ILE:N	2.82	0.42
1:A:227:VAL:CG2	1:A:228:ILE:N	2.82	0.42
1:B:214:ILE:HD12	1:B:219:MET:HG3	2.02	0.42
1:A:46:HIS:HD1	1:A:74:SER:CB	2.31	0.42
1:A:212:GLU:CD	1:A:252:VAL:HG11	2.39	0.42
1:B:161:LEU:HA	1:B:161:LEU:HD12	1.93	0.42
1:B:164:ARG:HG2	1:B:227:VAL:HB	2.02	0.42
1:A:176:TRP:CZ3	1:A:219:MET:HE3	2.55	0.42
1:A:262:VAL:CG1	1:A:263:THR:N	2.82	0.42
1:A:164:ARG:HG2	1:A:227:VAL:HB	2.02	0.42
1:B:135:ASP:O	1:B:136:GLN:C	2.59	0.42
1:B:93:TYR:HB3	1:B:100:PHE:HB3	2.01	0.41
1:A:93:TYR:CE1	1:A:256:PRO:HD3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:MET:O	1:B:177:LEU:HD23	2.20	0.41
1:B:252:VAL:C	1:B:254:GLY:H	2.24	0.41
1:B:304:LYS:O	1:B:308:LYS:HB2	2.21	0.41
1:B:88:ILE:HG22	1:B:93:TYR:HB2	2.02	0.41
1:A:88:ILE:HG22	1:A:93:TYR:HB2	2.02	0.41
1:B:114:GLN:HE21	1:B:114:GLN:H	1.68	0.41
1:A:165:MET:O	1:A:177:LEU:HD23	2.20	0.41
1:A:139:MET:HG2	1:A:295:ALA:HB3	2.03	0.41
1:A:93:TYR:HB3	1:A:100:PHE:HB3	2.02	0.41
1:B:262:VAL:HG12	1:B:263:THR:N	2.35	0.41
1:B:355:ASP:OD2	1:B:360:VAL:HG21	2.20	0.41
1:A:318:GLN:NE2	1:B:20:MET:O	2.54	0.41
1:A:375:THR:HG23	1:A:380:HIS:CD2	2.56	0.41
1:A:304:LYS:O	1:A:308:LYS:HB2	2.21	0.41
1:B:52:ASN:HA	1:B:288:TYR:HH	1.86	0.41
1:B:215:THR:HG23	1:B:218:ALA:CB	2.50	0.41
1:B:375:THR:HG23	1:B:380:HIS:CD2	2.55	0.41
1:A:214:ILE:HD12	1:A:219:MET:HG3	2.02	0.41
1:A:311:LEU:HB3	1:A:346:LEU:HD21	2.02	0.41
1:B:139:MET:HG2	1:B:295:ALA:HB3	2.02	0.41
1:A:209:GLN:HA	1:A:251:PHE:O	2.21	0.40
1:B:371:ILE:HD13	1:B:371:ILE:H	1.86	0.40
1:B:161:LEU:HD23	1:B:183:THR:CG2	2.52	0.40
1:B:311:LEU:HB3	1:B:346:LEU:HD21	2.04	0.40
1:B:72:ILE:HD11	1:B:80:TYR:HH	1.86	0.40
1:B:257:GLN:O	1:B:258:GLY:C	2.59	0.40
1:A:214:ILE:HD12	1:A:219:MET:HG2	2.04	0.40
1:A:219:MET:HE1	1:A:246:GLN:NE2	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	364/396 (92%)	310 (85%)	40 (11%)	14 (4%)	4	35
1	B	364/396 (92%)	312 (86%)	36 (10%)	16 (4%)	3	30
All	All	728/792 (92%)	622 (85%)	76 (10%)	30 (4%)	3	33

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ALA
1	A	193	ASN
1	A	196	VAL
1	B	18	ALA
1	B	193	ASN
1	B	196	VAL
1	A	344	ARG
1	B	135	ASP
1	B	258	GLY
1	B	344	ARG
1	A	115	SER
1	A	135	ASP
1	A	342	THR
1	B	63	THR
1	B	115	SER
1	B	342	THR
1	A	63	THR
1	A	195	ALA
1	A	260	ALA
1	A	330	SER
1	B	195	ALA
1	B	330	SER
1	A	131	VAL
1	B	131	VAL
1	B	253	ILE
1	B	256	PRO
1	A	258	GLY
1	A	194	GLY
1	B	194	GLY
1	B	252	VAL

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	307/332 (92%)	276 (90%)	31 (10%)	9	40
1	B	307/332 (92%)	275 (90%)	32 (10%)	9	39
All	All	614/664 (92%)	551 (90%)	63 (10%)	9	40

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	PHE
1	A	63	THR
1	A	67	LEU
1	A	68	LEU
1	A	100	PHE
1	A	114	GLN
1	A	139	MET
1	A	157	LEU
1	A	164	ARG
1	A	186	THR
1	A	191	GLN
1	A	203	THR
1	A	206	ILE
1	A	214	ILE
1	A	215	THR
1	A	227	VAL
1	A	239	GLU
1	A	251	PHE
1	A	265	ARG
1	A	267	ILE
1	A	275	TRP
1	A	289	THR
1	A	315	VAL
1	A	331	ILE
1	A	342	THR
1	A	343	GLU
1	A	344	ARG
1	A	369	LYS
1	A	371	ILE
1	A	375	THR
1	A	393	LYS

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Mol	Chain	Res	Type
1	B	19	PHE
1	B	63	THR
1	B	67	LEU
1	B	68	LEU
1	B	100	PHE
1	B	114	GLN
1	B	139	MET
1	B	157	LEU
1	B	164	ARG
1	B	186	THR
1	B	190	VAL
1	B	191	GLN
1	B	203	THR
1	B	206	ILE
1	B	214	ILE
1	B	227	VAL
1	B	239	GLU
1	B	251	PHE
1	B	257	GLN
1	B	265	ARG
1	B	267	ILE
1	B	275	TRP
1	B	289	THR
1	B	315	VAL
1	B	331	ILE
1	B	342	THR
1	B	343	GLU
1	B	344	ARG
1	B	369	LYS
1	B	371	ILE
1	B	375	THR
1	B	393	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	113	GLN
1	A	114	GLN
1	A	136	GLN
1	A	159	HIS
1	A	162	ASN

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Mol	Chain	Res	Type
1	A	191	GLN
1	A	209	GLN
1	A	210	HIS
1	A	278	HIS
1	A	318	GLN
1	A	353	ASN
1	A	380	HIS
1	B	52	ASN
1	B	113	GLN
1	B	114	GLN
1	B	136	GLN
1	B	159	HIS
1	B	162	ASN
1	B	184	GLN
1	B	191	GLN
1	B	209	GLN
1	B	210	HIS
1	B	278	HIS
1	B	318	GLN
1	B	353	ASN
1	B	380	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 are monoatomic - leaving 4 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	$\# Z > 2$
2	LIS	A	1398	4	6,9,9	1.12	0	6,12,12	0.92	1 (16%)
3	PO4	A	1400	4	4,4,4	0.68	0	6,6,6	0.27	0
6	ATP	B	1397	4	24,33,33	2.42	6 (25%)	31,52,52	3.26	11 (35%)
3	PO4	B	1399	-	4,4,4	0.75	0	6,6,6	0.37	0

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	LIS	A	1398	4	-	0/4/10/10	0/0/1/1
3	PO4	A	1400	4	-	0/0/0/0	0/0/0/0
6	ATP	B	1397	4	-	0/18/38/38	0/3/3/3
3	PO4	B	1399	-	-	0/0/0/0	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1397	ATP	C3'-C4'	2.06	1.58	1.53
6	B	1397	ATP	O4'-C4'	3.48	1.53	1.45
6	B	1397	ATP	PA-O5'	3.48	1.74	1.59
6	B	1397	ATP	C5'-C4'	4.12	1.65	1.51
6	B	1397	ATP	C2-N3	5.02	1.41	1.32
6	B	1397	ATP	O4'-C1'	7.35	1.50	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	1397	ATP	N3-C2-N1	-11.53	120.07	128.89
6	B	1397	ATP	O5'-PA-O1A	-2.08	101.53	109.62
6	B	1397	ATP	C2'-C3'-C4'	2.04	106.80	102.61
6	B	1397	ATP	C2'-C1'-N9	2.06	117.44	114.29
2	A	1398	LIS	CB-CA-N	2.20	115.37	109.91
6	B	1397	ATP	O3G-PG-O3B	2.23	115.23	105.09
6	B	1397	ATP	O2B-PB-O3B	2.90	118.27	105.09
6	B	1397	ATP	C4-C5-N7	3.02	112.25	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	1397	ATP	PA-O3A-PB	3.14	141.56	132.73
6	B	1397	ATP	O5'-C5'-C4'	5.26	128.50	109.12
6	B	1397	ATP	O3A-PA-O5'	6.81	121.01	102.94
6	B	1397	ATP	O4'-C1'-N9	7.67	124.16	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1398	LIS	1	0
3	A	1400	PO4	1	0
6	B	1397	ATP	13	0
3	B	1399	PO4	2	0

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.