



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 02:34 PM GMT

PDB ID : 1Q0T
Title : Ternary Structure of T4DAM with AdoHcy and DNA
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Deposited on : 2003-07-17
Resolution : 3.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

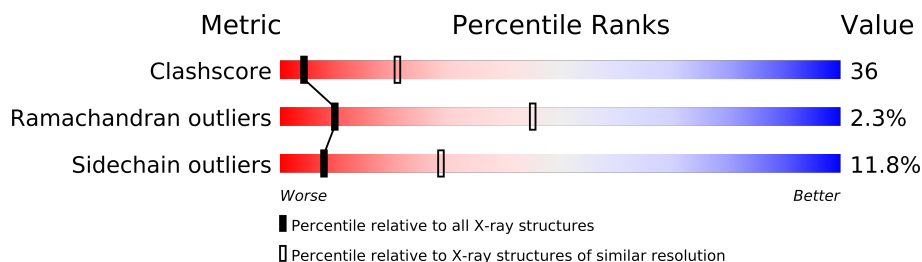
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : NOT EXECUTED
EDS : NOT EXECUTED
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.10 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	79885	1078 (3.16-3.04)
Ramachandran outliers	78287	1044 (3.16-3.04)
Sidechain outliers	78261	1044 (3.16-3.04)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	12	
1	D	12	
2	A	259	
2	B	259	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4538 atoms, of which 0 are hydrogen and 0 are deuterium.

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 DNA chain called 5'-D(*AP*CP*AP*GP*GP*AP*TP*CP*CP*TP*GP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	12	Total	C	N	O	P	0	0	0
			243	117	45	70	11			
1	D	12	Total	C	N	O	P	0	0	0
			243	117	45	70	11			

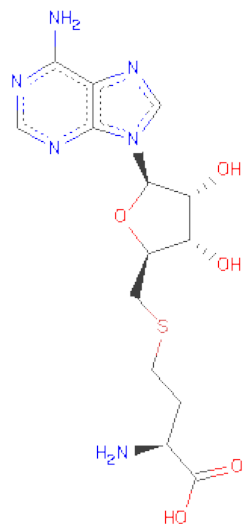
- Molecule 2 is a protein called DNA adenine methylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	241	Total	C	N	O	S	0	0	0
			1999	1293	335	366	5			
2	B	241	Total	C	N	O	S	0	0	0
			1999	1293	335	366	5			

- Molecule 3 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	I	0	0
			1	1		
3	C	1	Total	I	0	0
			1	1		

- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: C₁₄H₂₀N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
4	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

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.

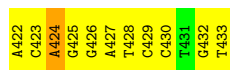
- Molecule 1: 5'-D(*AP*CP*AP*GP*GP*AP*TP*CP*CP*TP*GP*T)-3'

Chain C: 



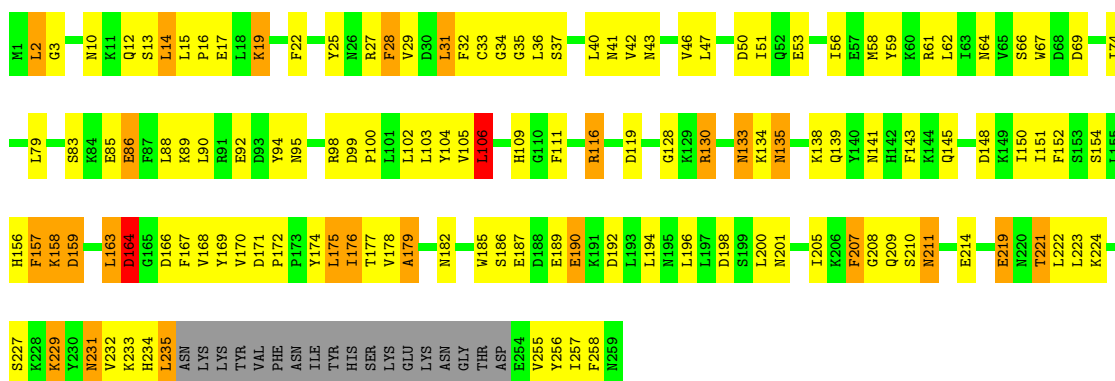
- Molecule 1: 5'-D(*AP*CP*AP*GP*GP*AP*TP*CP*CP*TP*GP*T)-3'

Chain D: 



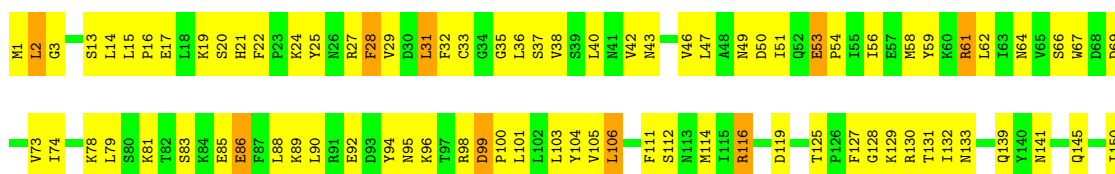
- Molecule 2: DNA adenine methylase

Chain A: 



- Molecule 2: DNA adenine methylase

Chain B: 



N220	L223	L224	E225	K226	S227	K228	K229	K230	N231	V232	V233	E234	L235	ASN	LYS	LYS	LYS	THR	VAL	PHE	ASN	ASN	ILE	THR	HIS	SER	LYS	GLY	GLU	LYS	ASN	ASN	GLY	THR	THR	ASP	E254	V255	V256	F258	N259										
L151	F152	S153	H156	F157	K158	D159	L163	D164	G165	D166	F167	V168	V169	V170	D171	D172	F173	V174	L175	L176	L177	L178	L181	L182	L183	F184	S186	E187	D188	E189	E190	L193	L194	N195	L196	L200	N201	D202	L203	G204	L205	L206	F207	G208	Q209	S210	N211	V212	L213	E214	F216

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.70Å 109.70Å 73.60Å 90.00° 104.20° 90.00°	Depositor
Resolution (Å)	20.00 – 3.10	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.10)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.238 , 0.293	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4538	wwPDB-VP
Average B, all atoms (Å ²)	40.0	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: SAH, IOD

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	C	0.74	0/272	1.45	5/418 (1.2%)
1	D	0.70	0/272	0.87	0/418
2	A	0.45	0/2046	0.65	0/2756
2	B	0.42	0/2046	0.63	0/2756
All	All	0.48	0/4636	0.74	5/6348 (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	5
1	D	0	1
All	All	0	6

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	405	DG	N9-C1'-C2'	7.85	127.52	112.60
1	C	408	DT	O4'-C1'-N1	6.32	112.42	108.00
1	C	406	DG	N9-C1'-C2'	6.11	124.21	112.60
1	C	408	DT	O3'-P-O5'	5.91	115.23	104.00
1	C	410	DC	N1-C1'-C2'	5.47	122.99	112.60

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	404	DA	Sidechain
1	C	406	DG	Sidechain
1	C	407	DA	Sidechain
1	C	409	DC	Sidechain
1	C	410	DC	Sidechain
1	D	424	DA	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	243	0	137	30	0
1	D	243	0	137	15	0
2	A	1999	0	1997	141	0
2	B	1999	0	1997	137	0
3	C	1	0	0	2	0
3	D	1	0	0	2	0
4	A	26	0	19	4	0
4	B	26	0	19	2	0
All	All	4538	0	4306	320	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 36.

All (320) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:423:DC:H5	3:D:1:IOD:I	1.94	1.21
2:B:83:SER:HB2	2:B:86:GLU:HG3	1.28	1.12
2:A:83:SER:HB2	2:A:86:GLU:HG3	1.35	1.08
1:C:403:DC:H5	3:C:2:IOD:I	2.08	1.07
1:D:423:DC:C5	3:D:1:IOD:I	2.78	1.06
2:A:50:ASP:HB3	2:A:56:ILE:HD11	1.40	1.04
2:A:15:LEU:HG	2:A:19:LYS:HD2	1.38	1.02
2:B:50:ASP:HB3	2:B:56:ILE:HD11	1.48	0.96
2:A:83:SER:HB2	2:A:86:GLU:CG	1.96	0.95
1:C:402:DA:H2''	1:C:403:DC:O5'	1.66	0.95

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:83:SER:HB2	2:B:86:GLU:CG	1.96	0.94
2:B:21:HIS:NE2	2:B:233:LYS:HE2	1.83	0.94
2:B:61:ARG:HH11	2:B:61:ARG:HG3	1.32	0.93
2:A:135:ASN:HD22	2:A:135:ASN:H	1.08	0.93
2:A:170:VAL:HG22	2:A:172:PRO:HD3	1.51	0.91
2:A:53:GLU:HG2	2:A:154:SER:CB	2.01	0.91
2:B:170:VAL:HG22	2:B:172:PRO:HD3	1.51	0.91
1:C:404:DA:H2''	1:C:405:DG:H5'	1.56	0.88
2:A:53:GLU:HG2	2:A:154:SER:HB2	1.58	0.85
2:A:141:ASN:O	2:A:145:GLN:HG2	1.78	0.84
1:C:403:DC:C5	3:C:2:IOD:I	2.99	0.82
2:A:14:LEU:HD12	2:A:14:LEU:H	1.44	0.81
2:A:135:ASN:N	2:A:135:ASN:HD22	1.80	0.79
2:B:19:LYS:HA	2:B:22:PHE:CD2	2.20	0.77
2:A:25:TYR:CE1	2:A:28:PHE:HB2	2.21	0.76
2:B:25:TYR:CE1	2:B:28:PHE:HB2	2.21	0.76
2:B:226:TRP:HA	2:B:229:LYS:HE3	1.67	0.76
2:A:232:VAL:HG12	2:A:257:ILE:HG12	1.68	0.76
2:B:232:VAL:HG12	2:B:257:ILE:HG12	1.67	0.76
2:B:141:ASN:O	2:B:145:GLN:HG2	1.86	0.75
2:A:135:ASN:H	2:A:135:ASN:ND2	1.84	0.74
2:B:169:TYR:HE2	2:B:171:ASP:HB2	1.54	0.72
1:C:408:DT:H2''	1:C:409:DC:O5'	1.88	0.72
2:A:85:GLU:OE2	2:A:85:GLU:HA	1.88	0.72
2:B:51:ILE:HB	4:B:301:SAH:C2	2.20	0.72
2:A:15:LEU:CG	2:A:19:LYS:HD2	2.20	0.71
2:A:164:ASP:OD2	2:A:164:ASP:N	2.23	0.71
2:A:17:GLU:OE2	2:A:235:LEU:HB3	1.91	0.71
2:B:21:HIS:CD2	2:B:233:LYS:HE2	2.27	0.70
2:B:15:LEU:HB3	2:B:16:PRO:HD3	1.75	0.69
2:A:33:CYS:HB2	2:A:36:LEU:HD23	1.74	0.69
1:C:412:DG:H2''	1:C:413:DT:C5'	2.24	0.68
1:C:407:DA:H2''	1:C:408:DT:C5'	2.23	0.68
2:B:14:LEU:HD12	2:B:14:LEU:H	1.57	0.68
2:B:61:ARG:NH1	2:B:61:ARG:HG3	2.03	0.68
2:A:35:GLY:O	2:A:36:LEU:HB2	1.95	0.67
2:B:33:CYS:HB2	2:B:36:LEU:HD23	1.76	0.67
1:C:403:DC:H2''	1:C:404:DA:O5'	1.92	0.67
2:B:209:GLN:HE21	2:B:210:SER:N	1.93	0.67
2:B:186:SER:OG	2:B:189:GLU:HG2	1.95	0.66
2:B:226:TRP:O	2:B:229:LYS:HG3	1.94	0.66
1:C:411:DT:OP2	2:A:133:ASN:HB2	1.94	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:32:PHE:CE1	2:B:170:VAL:HG23	2.31	0.66
2:A:32:PHE:CE1	2:A:170:VAL:HG23	2.31	0.66
1:D:425:DG:H1'	1:D:426:DG:C8	2.32	0.65
1:D:423:DC:H2''	1:D:424:DA:C8	2.32	0.65
2:B:29:VAL:CG1	2:B:168:VAL:HG22	2.28	0.64
2:A:29:VAL:CG1	2:A:168:VAL:HG22	2.28	0.64
2:A:14:LEU:HD12	2:A:14:LEU:N	2.13	0.63
2:A:190:GLU:O	2:A:194:LEU:HG	1.98	0.63
2:A:209:GLN:NE2	2:A:210:SER:O	2.31	0.63
2:A:169:TYR:HE2	2:A:171:ASP:HB2	1.63	0.63
2:A:135:ASN:N	2:A:135:ASN:ND2	2.41	0.63
2:B:94:TYR:CE1	2:B:103:LEU:HB2	2.34	0.63
2:A:25:TYR:HE1	2:A:28:PHE:HB2	1.63	0.62
2:A:15:LEU:HB3	2:A:16:PRO:HD3	1.80	0.62
2:B:235:LEU:H	2:B:235:LEU:HD12	1.65	0.62
2:B:163:LEU:O	2:B:164:ASP:O	2.18	0.62
2:B:209:GLN:HB3	2:B:257:ILE:O	2.00	0.61
2:A:158:LYS:HG2	2:A:159:ASP:N	2.15	0.61
1:C:407:DA:H2''	1:C:408:DT:H5'	1.83	0.60
2:A:74:ILE:HG23	2:A:79:LEU:HB2	1.83	0.60
2:A:163:LEU:O	2:A:164:ASP:O	2.20	0.60
1:C:406:DG:H2''	1:C:407:DA:O5'	2.02	0.60
2:A:61:ARG:HG3	2:A:61:ARG:HH11	1.66	0.59
2:B:227:SER:HB3	2:B:257:ILE:HD13	1.83	0.59
2:B:40:LEU:HD23	2:B:150:ILE:HD11	1.85	0.59
2:B:90:LEU:O	2:B:90:LEU:HD12	2.02	0.59
2:A:186:SER:OG	2:A:189:GLU:HG2	2.02	0.59
2:A:83:SER:HB2	2:A:86:GLU:HG2	1.84	0.59
2:B:190:GLU:O	2:B:194:LEU:HG	2.02	0.59
1:C:404:DA:H2'	1:C:405:DG:C8	2.39	0.58
2:A:209:GLN:HE21	2:A:210:SER:N	2.01	0.58
2:B:35:GLY:O	2:B:36:LEU:HB2	2.02	0.58
2:B:47:LEU:HD12	2:B:151:ILE:O	2.02	0.58
1:D:429:DC:H2''	1:D:430:DC:H5'	1.86	0.58
2:B:74:ILE:HG23	2:B:79:LEU:HB2	1.85	0.58
2:B:127:PHE:CE2	2:B:129:LYS:HA	2.38	0.58
2:A:19:LYS:HA	2:A:22:PHE:CD2	2.39	0.57
2:A:94:TYR:O	2:A:98:ARG:HD3	2.05	0.57
1:D:433:DT:H5'	1:D:433:DT:H6	1.68	0.57
2:B:169:TYR:CE2	2:B:171:ASP:HB2	2.39	0.57
2:A:169:TYR:C	2:A:169:TYR:CD2	2.77	0.56
2:B:17:GLU:OE2	2:B:235:LEU:HB3	2.05	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:3:GLY:HA2	2:A:37:SER:HA	1.87	0.56
2:B:3:GLY:HA2	2:B:37:SER:HA	1.87	0.56
2:B:58:MET:HA	2:B:100:PRO:HB3	1.87	0.56
1:C:412:DG:H2''	1:C:413:DT:H5'	1.87	0.56
2:A:58:MET:HA	2:A:100:PRO:HB3	1.86	0.56
1:C:404:DA:H2''	1:C:405:DG:C5'	2.34	0.56
2:A:50:ASP:CB	2:A:56:ILE:HD11	2.25	0.56
2:B:170:VAL:CG2	2:B:172:PRO:HD3	2.29	0.56
2:B:209:GLN:NE2	2:B:210:SER:O	2.38	0.56
1:C:412:DG:H2''	1:C:413:DT:O5'	2.06	0.56
1:C:402:DA:H2''	1:C:403:DC:C5'	2.35	0.55
2:B:17:GLU:OE1	2:B:233:LYS:HE3	2.06	0.55
2:B:27:ARG:NH2	2:B:163:LEU:H	2.05	0.55
2:B:178:VAL:O	2:B:179:ALA:HB2	2.05	0.55
1:C:409:DC:H2''	1:C:410:DC:C5'	2.36	0.55
2:B:169:TYR:C	2:B:169:TYR:CD2	2.80	0.55
2:B:116:ARG:HG3	2:B:125:THR:OG1	2.07	0.55
2:B:175:LEU:HD12	2:B:190:GLU:HG3	1.88	0.55
2:A:175:LEU:HD12	2:A:190:GLU:HG3	1.90	0.54
2:B:214:GLU:HA	2:B:219:GLU:HA	1.88	0.54
2:A:2:LEU:HD22	2:A:139:GLN:HG2	1.87	0.54
1:C:409:DC:H2''	1:C:410:DC:H5'	1.89	0.54
2:A:134:LYS:O	2:A:138:LYS:HG3	2.06	0.54
2:B:158:LYS:HE2	2:B:159:ASP:OD1	2.07	0.54
2:A:14:LEU:CD1	2:A:14:LEU:H	2.19	0.54
1:D:427:DA:H1'	1:D:428:DT:H5''	1.88	0.54
2:A:66:SER:O	2:A:69:ASP:HB2	2.08	0.54
2:A:62:LEU:HD11	2:A:104:TYR:CE1	2.42	0.54
2:A:28:PHE:HB3	2:A:46:VAL:HG22	1.90	0.54
2:A:10:ASN:CG	2:A:12:GLN:HG3	2.29	0.54
2:A:214:GLU:HA	2:A:219:GLU:HA	1.89	0.54
2:A:111:PHE:HA	2:A:128:GLY:HA2	1.91	0.53
2:B:25:TYR:HE1	2:B:28:PHE:HB2	1.70	0.53
2:A:31:LEU:CD2	2:A:157:PHE:HB2	2.38	0.53
2:A:50:ASP:OD2	4:A:300:SAH:O2'	2.26	0.53
2:A:99:ASP:HB3	2:A:102:LEU:HD12	1.90	0.53
1:D:423:DC:H5'	1:D:423:DC:H6	1.74	0.53
2:B:50:ASP:CB	2:B:56:ILE:HD11	2.31	0.53
2:B:235:LEU:HD12	2:B:235:LEU:N	2.24	0.52
2:B:2:LEU:HD22	2:B:139:GLN:HG2	1.90	0.52
2:A:27:ARG:NH2	2:A:163:LEU:H	2.07	0.52
2:A:13:SER:O	2:A:16:PRO:HD2	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:407:DA:H2''	1:C:408:DT:O5'	2.10	0.52
2:B:213:LEU:CD1	2:B:255:VAL:HG11	2.40	0.52
1:C:406:DG:H2''	1:C:407:DA:C5'	2.39	0.52
1:C:408:DT:H4'	1:C:409:DC:OP1	2.10	0.52
2:A:233:LYS:O	2:A:255:VAL:HA	2.10	0.52
2:B:209:GLN:NE2	2:B:210:SER:N	2.57	0.52
2:B:92:GLU:O	2:B:96:LYS:HG2	2.10	0.51
1:D:425:DG:H4'	1:D:426:DG:OP1	2.10	0.51
2:B:53:GLU:HB2	2:B:54:PRO:HD3	1.90	0.51
2:B:233:LYS:O	2:B:255:VAL:HA	2.11	0.51
2:B:29:VAL:HG11	2:B:168:VAL:HG22	1.91	0.51
2:B:223:LEU:HD23	2:B:223:LEU:O	2.10	0.51
2:A:32:PHE:CD1	2:A:170:VAL:HG23	2.45	0.51
2:A:223:LEU:O	2:A:223:LEU:HD23	2.10	0.51
2:A:150:ILE:HG21	2:A:152:PHE:CE2	2.46	0.51
2:A:116:ARG:HG3	2:A:116:ARG:NH1	2.24	0.51
2:B:166:ASP:N	2:B:166:ASP:OD1	2.43	0.51
2:A:40:LEU:HD23	2:A:150:ILE:HD11	1.92	0.51
2:B:111:PHE:O	2:B:112:SER:HB2	2.12	0.51
2:B:167:PHE:HZ	2:B:256:TYR:OH	1.94	0.50
1:C:406:DG:H2'	1:C:407:DA:C8	2.45	0.50
2:B:14:LEU:HD12	2:B:14:LEU:N	2.25	0.50
2:B:66:SER:O	2:B:69:ASP:HB2	2.11	0.50
1:D:432:DG:H1'	1:D:433:DT:H5''	1.93	0.50
2:B:156:HIS:CD2	2:B:158:LYS:H	2.29	0.50
2:A:99:ASP:C	2:A:99:ASP:OD1	2.48	0.50
2:A:227:SER:HB3	2:A:257:ILE:HD13	1.93	0.50
1:C:406:DG:OP1	2:B:133:ASN:HB2	2.11	0.50
2:A:235:LEU:HD12	2:A:235:LEU:H	1.76	0.50
2:A:167:PHE:HZ	2:A:256:TYR:OH	1.94	0.50
2:B:38:VAL:O	2:B:42:VAL:HG13	2.11	0.50
2:A:95:ASN:O	2:A:98:ARG:NE	2.43	0.50
2:A:178:VAL:O	2:A:179:ALA:HB2	2.11	0.50
2:B:176:ILE:HG13	2:B:176:ILE:O	2.11	0.50
2:A:156:HIS:CD2	2:A:158:LYS:H	2.30	0.50
2:B:112:SER:HA	2:B:130:ARG:NH1	2.26	0.50
2:A:169:TYR:CE2	2:A:171:ASP:HB2	2.44	0.50
1:D:422:DA:H2''	1:D:423:DC:H5'	1.93	0.49
2:B:14:LEU:O	2:B:15:LEU:C	2.50	0.49
2:B:33:CYS:HG	2:B:49:ASN:C	2.16	0.49
2:A:179:ALA:N	2:A:182:ASN:HD22	2.10	0.49
2:A:207:PHE:CD2	2:A:207:PHE:C	2.85	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:196:LEU:O	2:B:196:LEU:HD12	2.13	0.49
2:A:233:LYS:HD3	2:A:234:HIS:H	1.78	0.49
2:B:207:PHE:C	2:B:207:PHE:CD2	2.86	0.49
2:B:56:ILE:O	2:B:59:TYR:HB2	2.13	0.49
2:A:130:ARG:HG3	2:A:130:ARG:HH11	1.76	0.49
2:A:14:LEU:O	2:A:15:LEU:C	2.51	0.49
2:A:166:ASP:N	2:A:166:ASP:OD1	2.44	0.48
2:A:47:LEU:HD12	2:A:151:ILE:O	2.13	0.48
2:A:67:TRP:CE3	2:A:67:TRP:HA	2.48	0.48
2:A:51:ILE:HB	4:A:300:SAH:C2	2.42	0.48
2:A:221:THR:HG22	2:A:222:LEU:N	2.29	0.48
2:B:29:VAL:HG12	2:B:167:PHE:O	2.13	0.48
2:B:99:ASP:OD1	2:B:99:ASP:C	2.52	0.48
2:A:85:GLU:OE2	2:A:88:LEU:HD12	2.13	0.48
2:A:85:GLU:OE2	2:A:85:GLU:CA	2.59	0.48
2:B:83:SER:HB2	2:B:86:GLU:HG2	1.91	0.48
2:B:233:LYS:HD3	2:B:235:LEU:HG	1.94	0.48
2:B:177:THR:O	2:B:182:ASN:ND2	2.47	0.48
2:A:89:LYS:HA	2:A:92:GLU:HB2	1.96	0.48
2:B:131:THR:HG23	2:B:132:ILE:N	2.28	0.47
2:B:180:ASP:O	2:B:183:LYS:HG2	2.14	0.47
2:A:31:LEU:HB3	2:A:32:PHE:CD1	2.50	0.47
2:B:193:LEU:O	2:B:196:LEU:HB3	2.13	0.47
2:B:33:CYS:HB2	2:B:36:LEU:CD2	2.42	0.47
2:A:169:TYR:CD2	2:A:170:VAL:N	2.83	0.47
2:B:179:ALA:N	2:B:182:ASN:HD22	2.12	0.47
2:A:61:ARG:HG3	2:A:61:ARG:NH1	2.28	0.47
2:B:159:ASP:N	2:B:159:ASP:OD1	2.47	0.47
2:A:10:ASN:OD1	2:A:12:GLN:HG3	2.15	0.47
2:A:29:VAL:HG11	2:A:168:VAL:HG22	1.96	0.47
2:B:200:LEU:O	2:B:205:ILE:HB	2.15	0.47
1:D:422:DA:H2''	1:D:423:DC:C5'	2.44	0.47
2:B:32:PHE:CD1	2:B:170:VAL:HG23	2.50	0.46
1:C:410:DC:OP1	2:A:130:ARG:NH2	2.49	0.46
2:B:67:TRP:CE3	2:B:67:TRP:HA	2.51	0.46
2:B:195:ASN:HD22	2:B:195:ASN:N	2.13	0.46
2:B:95:ASN:O	2:B:98:ARG:NE	2.45	0.46
2:B:62:LEU:HD11	2:B:104:TYR:CE1	2.51	0.46
2:A:209:GLN:HB3	2:A:257:ILE:O	2.16	0.46
2:B:235:LEU:HD12	2:B:254:GLU:O	2.15	0.46
1:C:406:DG:H2'	1:C:407:DA:H8	1.80	0.46
2:B:201:ASN:HD22	2:B:259:ASN:HD22	1.65	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:208:GLY:CA	2:A:256:TYR:CE1	2.99	0.45
2:A:209:GLN:O	2:A:256:TYR:HA	2.17	0.45
2:A:33:CYS:HB2	2:A:36:LEU:CD2	2.43	0.45
2:B:40:LEU:HD23	2:B:150:ILE:CD1	2.45	0.45
2:A:105:VAL:O	2:A:109:HIS:HD2	2.00	0.45
2:A:35:GLY:N	4:A:300:SAH:O	2.47	0.45
2:B:169:TYR:CD2	2:B:170:VAL:N	2.85	0.45
2:B:170:VAL:HG12	2:B:208:GLY:O	2.17	0.45
2:A:256:TYR:CE2	2:A:258:PHE:HB3	2.52	0.45
2:A:56:ILE:O	2:A:59:TYR:HB2	2.17	0.45
2:A:27:ARG:NH2	2:A:166:ASP:OD2	2.35	0.45
2:A:79:LEU:HD11	2:A:90:LEU:HD22	1.99	0.45
1:C:409:DC:H2''	1:C:410:DC:O5'	2.17	0.44
2:B:73:VAL:HG11	2:B:105:VAL:HG21	1.98	0.44
2:B:33:CYS:SG	2:B:49:ASN:O	2.74	0.44
2:A:40:LEU:HD23	2:A:150:ILE:CD1	2.48	0.44
2:A:85:GLU:O	2:A:88:LEU:HB2	2.17	0.44
2:B:111:PHE:HA	2:B:128:GLY:HA2	1.99	0.44
2:A:35:GLY:O	2:A:36:LEU:CB	2.65	0.44
2:A:198:ASP:O	2:A:201:ASN:HB3	2.18	0.44
2:B:29:VAL:O	2:B:29:VAL:HG13	2.17	0.44
1:C:412:DG:C2'	1:C:413:DT:O5'	2.66	0.44
1:D:425:DG:H1'	1:D:426:DG:N7	2.33	0.44
2:B:156:HIS:O	2:B:158:LYS:N	2.51	0.44
2:B:61:ARG:NH1	2:B:61:ARG:CG	2.70	0.44
2:B:28:PHE:HB3	2:B:46:VAL:HG22	2.00	0.44
2:B:15:LEU:O	2:B:16:PRO:C	2.56	0.44
2:B:13:SER:O	2:B:16:PRO:HD2	2.18	0.44
2:B:220:ASN:OD1	2:B:223:LEU:HB2	2.17	0.44
2:B:101:LEU:O	2:B:105:VAL:HG23	2.18	0.44
2:A:174:TYR:CE1	2:A:211:ASN:HA	2.53	0.43
2:A:176:ILE:HG13	2:A:176:ILE:O	2.17	0.43
2:B:85:GLU:O	2:B:88:LEU:HB2	2.18	0.43
2:A:42:VAL:O	2:A:43:ASN:C	2.56	0.43
2:B:49:ASN:HB2	2:B:153:SER:OG	2.18	0.43
2:B:94:TYR:O	2:B:98:ARG:HD3	2.18	0.43
2:A:111:PHE:HA	2:A:128:GLY:CA	2.48	0.43
2:A:31:LEU:HD21	2:A:157:PHE:HB2	2.00	0.43
2:B:89:LYS:HA	2:B:92:GLU:HB2	2.01	0.43
1:C:409:DC:C2'	1:C:410:DC:O5'	2.67	0.43
2:B:187:GLU:O	2:B:190:GLU:HB2	2.18	0.43
2:A:170:VAL:CG2	2:A:172:PRO:HD3	2.35	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:33:CYS:O	2:A:34:GLY:C	2.56	0.43
2:A:15:LEU:O	2:A:16:PRO:C	2.56	0.42
2:B:94:TYR:CZ	2:B:103:LEU:HB2	2.54	0.42
2:A:31:LEU:HD23	2:A:157:PHE:HB2	2.01	0.42
2:A:62:LEU:HD23	2:A:62:LEU:HA	1.74	0.42
2:B:209:GLN:O	2:B:256:TYR:HA	2.19	0.42
2:B:112:SER:HB2	2:B:114:MET:HG2	2.01	0.42
2:B:203:ARG:C	2:B:205:ILE:H	2.23	0.42
2:A:231:ASN:N	2:A:231:ASN:OD1	2.52	0.42
2:A:156:HIS:O	2:A:158:LYS:N	2.52	0.42
2:A:42:VAL:O	2:A:42:VAL:HG23	2.19	0.42
1:D:429:DC:H1'	1:D:430:DC:H5''	2.00	0.42
2:B:29:VAL:HG13	2:B:168:VAL:HG13	2.02	0.42
2:B:150:ILE:HG21	2:B:152:PHE:CE2	2.55	0.42
2:B:17:GLU:OE1	2:B:233:LYS:CE	2.68	0.42
2:B:24:LYS:O	2:B:25:TYR:HB3	2.20	0.42
2:A:15:LEU:HD21	2:A:41:ASN:HB2	2.02	0.42
2:B:31:LEU:HB3	2:B:32:PHE:CD1	2.55	0.42
2:A:27:ARG:O	2:A:166:ASP:HB2	2.20	0.42
2:A:53:GLU:HG2	2:A:154:SER:OG	2.19	0.41
2:A:159:ASP:N	2:A:159:ASP:OD1	2.52	0.41
2:A:177:THR:O	2:A:182:ASN:ND2	2.53	0.41
2:B:209:GLN:HE21	2:B:209:GLN:C	2.24	0.41
2:B:156:HIS:HD2	2:B:158:LYS:H	1.67	0.41
2:A:200:LEU:O	2:A:205:ILE:HB	2.20	0.41
2:A:106:LEU:HA	2:A:106:LEU:HD23	1.82	0.41
2:A:224:LYS:O	2:A:227:SER:OG	2.38	0.41
2:A:229:LYS:CB	2:A:229:LYS:NZ	2.83	0.41
1:C:410:DC:H2''	1:C:411:DT:O5'	2.19	0.41
2:A:187:GLU:O	2:A:190:GLU:HB2	2.20	0.41
2:A:171:ASP:OD2	2:A:171:ASP:C	2.59	0.41
2:A:209:GLN:HE21	2:A:210:SER:C	2.22	0.41
1:C:407:DA:C2'	1:C:408:DT:O5'	2.68	0.41
2:A:235:LEU:HD12	2:A:235:LEU:N	2.35	0.41
2:A:116:ARG:HH11	2:A:116:ARG:HG3	1.84	0.41
2:A:105:VAL:O	2:A:109:HIS:CD2	2.74	0.41
2:A:29:VAL:HG13	2:A:168:VAL:HA	2.03	0.41
2:A:100:PRO:O	2:A:103:LEU:N	2.53	0.41
2:B:195:ASN:N	2:B:195:ASN:ND2	2.68	0.41
2:A:32:PHE:HD2	4:A:300:SAH:O4'	2.04	0.41
2:B:224:LYS:O	2:B:227:SER:OG	2.39	0.41
2:B:163:LEU:O	2:B:166:ASP:OD1	2.37	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:170:VAL:HG12	2:A:208:GLY:O	2.21	0.41
2:B:33:CYS:SG	2:B:50:ASP:HB2	2.61	0.41
2:B:32:PHE:HD2	4:B:301:SAH:O4'	2.03	0.41
2:B:15:LEU:HB3	2:B:16:PRO:CD	2.48	0.41
2:A:156:HIS:HD2	2:A:158:LYS:H	1.68	0.41
2:B:42:VAL:HG23	2:B:42:VAL:O	2.21	0.41
2:B:62:LEU:HD23	2:B:62:LEU:HA	1.83	0.41
2:B:17:GLU:OE1	2:B:233:LYS:NZ	2.51	0.40
2:B:14:LEU:CD1	2:B:14:LEU:H	2.31	0.40
2:A:233:LYS:HD3	2:A:234:HIS:N	2.36	0.40
1:D:432:DG:H2''	1:D:433:DT:C5'	2.50	0.40
2:A:62:LEU:HB3	2:A:143:PHE:CE2	2.56	0.40
2:A:196:LEU:O	2:A:196:LEU:HD12	2.21	0.40
2:B:53:GLU:CB	2:B:54:PRO:HD3	2.50	0.40
2:B:174:TYR:HB2	2:B:177:THR:HB	2.04	0.40
2:B:42:VAL:O	2:B:43:ASN:C	2.59	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	237/259 (92%)	194 (82%)	37 (16%)	6 (2%)	9	42
2	B	237/259 (92%)	196 (83%)	36 (15%)	5 (2%)	11	48
All	All	474/518 (92%)	390 (82%)	73 (15%)	11 (2%)	10	45

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	164	ASP
2	B	164	ASP
2	A	157	PHE
2	A	179	ALA

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Mol	Chain	Res	Type
2	B	157	PHE
2	B	179	ALA
2	A	221	THR
2	B	106	LEU
2	A	106	LEU
2	A	176	ILE
2	B	176	ILE

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	
2	A	225/242 (93%)	197 (88%)	28 (12%)	7	25
2	B	225/242 (93%)	200 (89%)	25 (11%)	9	33
All	All	450/484 (93%)	397 (88%)	53 (12%)	8	29

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	2	LEU
2	A	14	LEU
2	A	19	LYS
2	A	28	PHE
2	A	31	LEU
2	A	64	ASN
2	A	86	GLU
2	A	106	LEU
2	A	116	ARG
2	A	119	ASP
2	A	130	ARG
2	A	133	ASN
2	A	135	ASN
2	A	148	ASP
2	A	158	LYS
2	A	159	ASP
2	A	163	LEU
2	A	164	ASP

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Mol	Chain	Res	Type
2	A	175	LEU
2	A	185	TRP
2	A	190	GLU
2	A	192	ASP
2	A	207	PHE
2	A	211	ASN
2	A	219	GLU
2	A	229	LYS
2	A	231	ASN
2	A	235	LEU
2	B	1	MET
2	B	2	LEU
2	B	20	SER
2	B	28	PHE
2	B	31	LEU
2	B	53	GLU
2	B	61	ARG
2	B	64	ASN
2	B	78	LYS
2	B	81	LYS
2	B	86	GLU
2	B	99	ASP
2	B	106	LEU
2	B	116	ARG
2	B	119	ASP
2	B	159	ASP
2	B	163	LEU
2	B	164	ASP
2	B	175	LEU
2	B	185	TRP
2	B	190	GLU
2	B	207	PHE
2	B	211	ASN
2	B	231	ASN
2	B	235	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	12	GLN
2	A	43	ASN
2	A	109	HIS

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Mol	Chain	Res	Type
2	A	135	ASN
2	A	139	GLN
2	A	141	ASN
2	A	156	HIS
2	A	182	ASN
2	A	201	ASN
2	A	209	GLN
2	B	12	GLN
2	B	43	ASN
2	B	109	HIS
2	B	139	GLN
2	B	141	ASN
2	B	156	HIS
2	B	182	ASN
2	B	201	ASN
2	B	209	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SAH	A	300	-	28,28,28	0.76	1 (3%)	40,40,40	1.11	3 (7%)
4	SAH	B	301	-	28,28,28	0.80	1 (3%)	40,40,40	1.12	2 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SAH	A	300	-	-	0/15/31/31	0/1/3/3
4	SAH	B	301	-	-	0/15/31/31	0/1/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301	SAH	O-C	2.84	1.31	1.22
4	A	300	SAH	O-C	2.53	1.30	1.22

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	SAH	C-CA-N	-4.38	102.11	109.36
4	A	300	SAH	C5'-SD-CG	3.76	113.64	102.42
4	B	301	SAH	C4'-C5'-SD	-2.97	104.38	113.53
4	A	300	SAH	OXT-C-O	-2.61	118.17	124.07
4	A	300	SAH	OXT-C-CA	2.14	121.69	116.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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.