



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:46 PM GMT

PDB ID : 1H5T
Title : THYMIDYLYLTRANSFERASE COMPLEXED WITH THYMIDYLYLDIPHOSPHATE-GLUCOSE
Authors : Rosano, C.; Zuccotti, S.; Bolognesi, M.
Deposited on : 2001-05-25
Resolution : 1.90 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

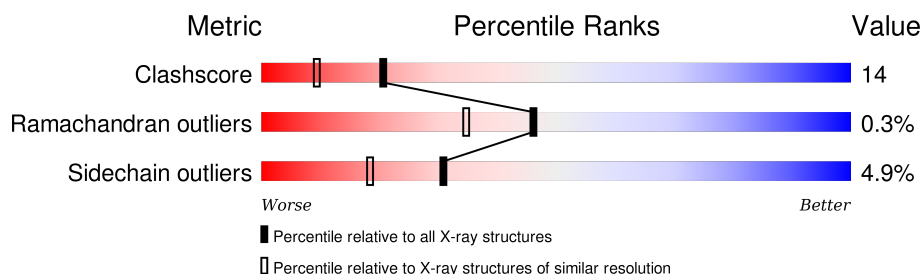
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)

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	293	
1	C	293	
1	D	293	
2	B	293	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9839 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 GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERAS E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	290	Total	C	N	O	S	25	0	0
			2277	1459	378	429	11			
1	C	290	Total	C	N	O	S	16	0	0
			2269	1455	378	425	11			
1	D	290	Total	C	N	O	S	10	0	0
			2273	1457	378	427	11			

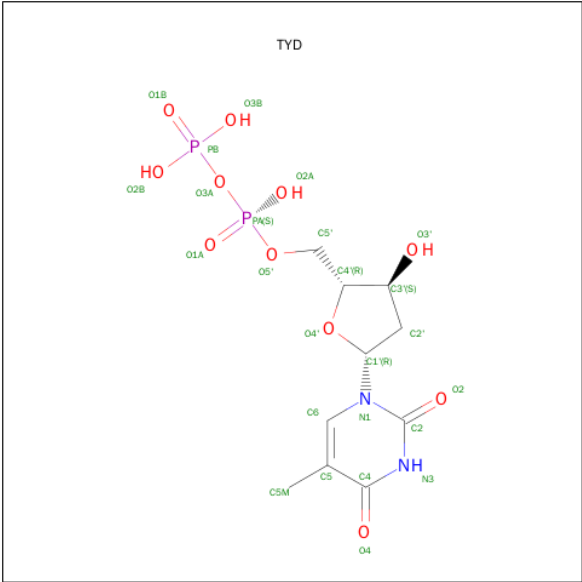
- Molecule 2 is a protein called GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERAS E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	290	Total	C	N	O	S	18	0	0
			2270	1455	378	428	9			

There are 2 discrepancies between the modelled and reference sequences:

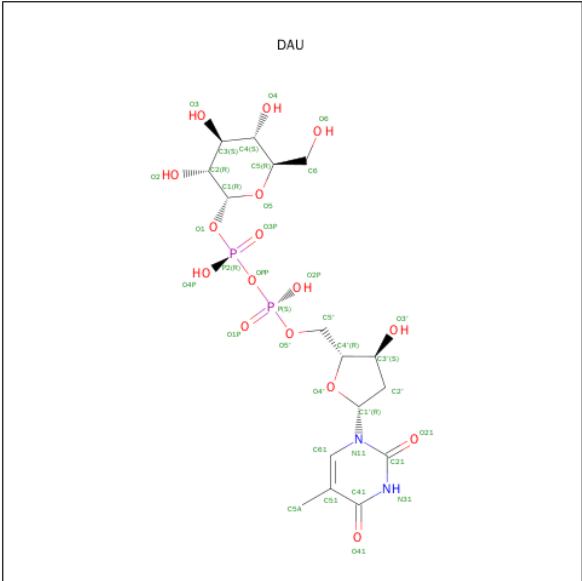
Chain	Residue	Modelled	Actual	Comment	Reference
B	217	LEU	MET	CONFLICT	UNP P37744
B	288	GLN	MET	CONFLICT	UNP P37744

- Molecule 3 is THYMIDINE-5'-DIPHOSPHATE (three-letter code: TYD) (formula: $C_{10}H_{16}N_2O_{11}P_2$).



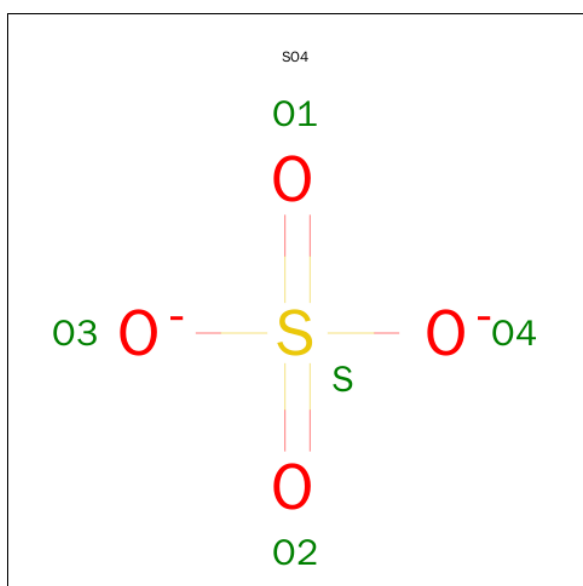
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
3	B	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
3	C	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
3	D	1	Total	C	N	O	P	0	0
			25	10	2	11	2		

- Molecule 4 is 2'DEOXY-THYMIDINE-5'-DIPHOSPHO-ALPHA-D-GLUCOSE (three-letter code: DAU) (formula: C₁₆H₂₆N₂O₁₆P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			36	16	2	16	2		
4	B	1	Total	C	N	O	P	0	0
			36	16	2	16	2		
4	C	1	Total	C	N	O	P	0	0
			36	16	2	16	2		
4	D	1	Total	C	N	O	P	0	0
			36	16	2	16	2		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

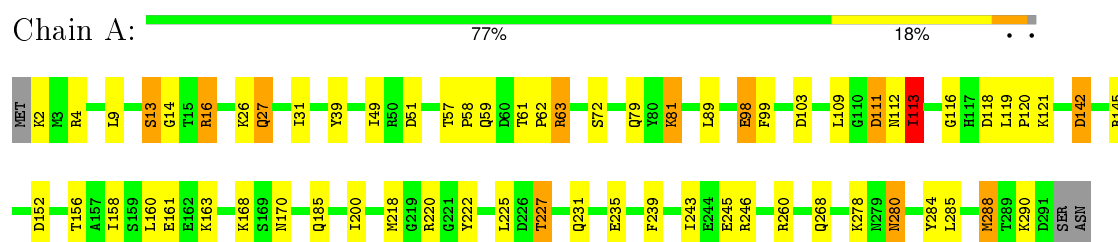
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	146	Total	O	0	0
			146	146		
6	B	99	Total	O	0	0
			99	99		
6	C	119	Total	O	0	0
			119	119		
6	D	137	Total	O	0	0
			137	137		

3 Residue-property plots

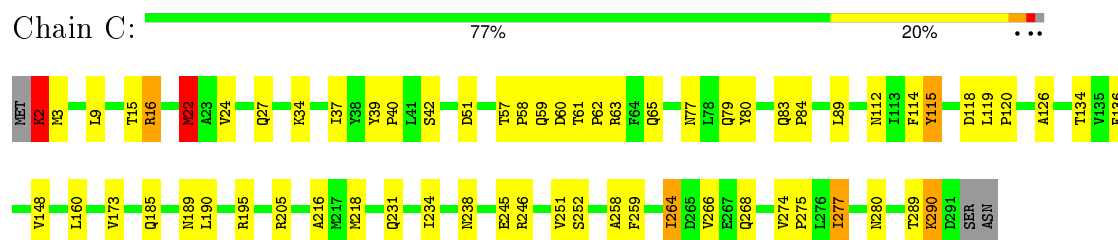
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

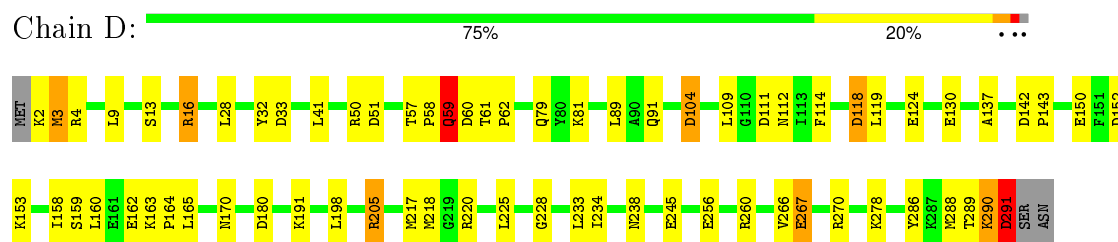
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



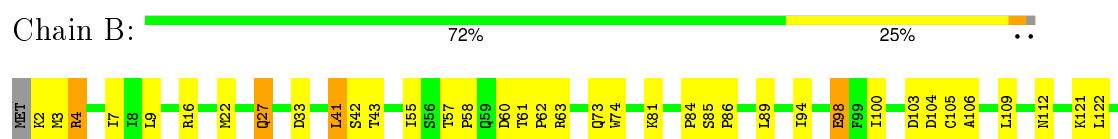
• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE

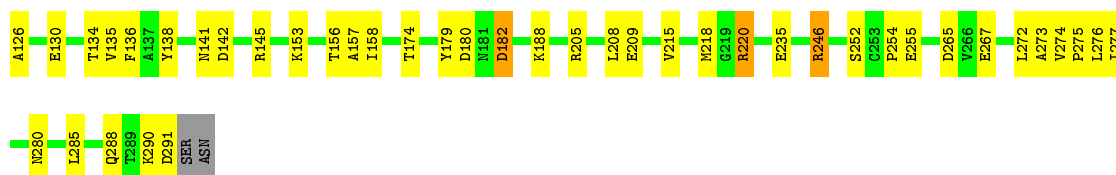


• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE



• Molecule 2: GLUCOSE-1-PHOSPHATE THYMIDYLTRANSFERASE





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	72.81Å 119.49Å 81.17Å 90.00° 112.67° 90.00°	Depositor
Resolution (Å)	12.00 – 1.90	Depositor
% Data completeness (in resolution range)	97.4 (12.00-1.90)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.174 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9839	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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: TYD, DAU, SO4

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	0.93	2/2324 (0.1%)	1.01	8/3145 (0.3%)
1	C	0.87	2/2316 (0.1%)	1.06	11/3134 (0.4%)
1	D	1.25	3/2320 (0.1%)	1.12	12/3140 (0.4%)
2	B	0.87	5/2317 (0.2%)	0.94	7/3138 (0.2%)
All	All	0.99	12/9277 (0.1%)	1.03	38/12557 (0.3%)

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	D	0	4
2	B	1	1
All	All	1	5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	59	GLN	CG-CD	42.29	2.48	1.51
1	D	290	LYS	CA-CB	15.68	1.88	1.53
1	A	245	GLU	CG-CD	-15.36	1.28	1.51
1	C	290	LYS	CB-CG	12.06	1.85	1.52
2	B	291	ASP	N-CA	-11.12	1.24	1.46
2	B	2	LYS	CD-CE	9.42	1.74	1.51
1	C	2	LYS	CD-CE	9.20	1.74	1.51
2	B	290	LYS	CB-CG	6.46	1.70	1.52
1	A	218	MET	SD-CE	-6.33	1.42	1.77
2	B	121	LYS	CG-CD	6.19	1.73	1.52
1	D	59	GLN	CB-CG	5.73	1.68	1.52
2	B	42	SER	CA-CB	5.04	1.60	1.52

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	59	GLN	CG-CD-NE2	-24.54	57.80	116.70
1	C	290	LYS	CB-CA-C	18.41	147.22	110.40
1	C	290	LYS	CA-CB-CG	-10.94	89.34	113.40
1	D	291	ASP	CB-CG-OD2	-9.01	110.19	118.30
1	C	290	LYS	CA-C-O	8.01	136.92	120.10
1	C	290	LYS	CA-C-N	-8.00	99.60	117.20
1	D	290	LYS	N-CA-CB	-7.61	96.91	110.60
1	A	16	ARG	CB-CG-CD	7.36	130.73	111.60
1	D	111	ASP	CB-CG-OD2	7.26	124.84	118.30
1	D	290	LYS	CB-CA-C	6.93	124.27	110.40
1	A	142	ASP	CB-CG-OD2	6.80	124.42	118.30
1	A	220	ARG	NE-CZ-NH2	-6.67	116.97	120.30
2	B	41	LEU	CB-CG-CD1	6.67	122.33	111.00
2	B	182	ASP	CB-CG-OD2	6.58	124.22	118.30
1	D	59	GLN	CB-CG-CD	6.51	128.52	111.60
1	D	60	ASP	CB-CG-OD2	6.35	124.01	118.30
2	B	60	ASP	CB-CG-OD2	6.06	123.75	118.30
1	C	148	VAL	CG1-CB-CG2	5.99	120.49	110.90
1	A	113	ILE	CG1-CB-CG2	5.96	124.52	111.40
2	B	265	ASP	CB-CG-OD2	5.96	123.67	118.30
1	C	2	LYS	CD-CE-NZ	5.96	125.41	111.70
2	B	246	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	D	33	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	51	ASP	CB-CG-OD2	5.78	123.50	118.30
1	C	60	ASP	CB-CG-OD2	5.71	123.44	118.30
1	D	142	ASP	CB-CG-OD2	5.64	123.38	118.30
1	A	111	ASP	CB-CG-OD2	5.63	123.36	118.30
1	C	277	ILE	CG1-CB-CG2	5.56	123.64	111.40
1	C	148	VAL	CA-CB-CG1	5.45	119.08	110.90
1	D	118	ASP	CB-CG-OD2	5.37	123.14	118.30
2	B	33	ASP	CB-CG-OD2	5.37	123.13	118.30
2	B	104	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	103	ASP	CB-CG-OD2	5.18	122.96	118.30
1	A	118	ASP	CB-CG-OD2	5.15	122.93	118.30
1	C	118	ASP	CB-CG-OD2	5.11	122.89	118.30
1	D	205	ARG	NE-CZ-NH1	-5.07	117.77	120.30
1	D	104	ASP	CB-CG-OD2	5.05	122.84	118.30
1	C	22	MET	CG-SD-CE	-5.02	92.17	100.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	174	THR	CB

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	106	ALA	Mainchain
1	D	137	ALA	Mainchain
1	D	267	GLU	Mainchain
1	D	291	ASP	Sidechain
1	D	59	GLN	Sidechain

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	2277	0	2284	65	0
1	C	2269	0	2275	70	0
1	D	2273	0	2280	67	0
2	B	2270	0	2271	71	0
3	A	25	0	13	0	0
3	B	25	0	13	0	0
3	C	25	0	13	0	0
3	D	25	0	13	0	0
4	A	36	0	16	2	0
4	B	36	0	16	1	0
4	C	36	0	16	2	0
4	D	36	0	16	1	0
5	D	5	0	0	1	0
6	A	146	0	0	12	1
6	B	99	0	0	8	1
6	C	119	0	0	5	1
6	D	137	0	0	14	1
All	All	9839	0	9226	265	2

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 (265) 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:C:112:ASN:ND2	1:C:114:PHE:CE1	1.74	1.55
1:D:217:MET:SD	1:D:217:MET:CE	2.01	1.46
2:B:142:ASP:OD2	2:B:145:ARG:NH1	1.67	1.27
1:A:63:ARG:CG	1:A:63:ARG:HH11	1.63	1.11
1:A:63:ARG:HH11	1:A:63:ARG:HG2	0.97	1.07
1:C:112:ASN:ND2	1:C:114:PHE:CZ	2.05	1.05
1:C:16:ARG:H	1:C:16:ARG:CD	1.70	1.04
1:A:113:ILE:CD1	1:A:225:LEU:HD12	1.85	1.04
1:C:16:ARG:H	1:C:16:ARG:HD2	1.18	1.02
1:D:130:GLU:OE2	6:D:2060:HOH:O	1.77	1.02
1:C:3:MET:HE3	1:C:51:ASP:HB2	1.42	1.00
2:B:235:GLU:OE1	6:B:2079:HOH:O	1.79	1.00
2:B:84:PRO:HG2	6:B:2034:HOH:O	1.62	1.00
2:B:98:GLU:H	2:B:98:GLU:CD	1.68	0.97
1:A:9:LEU:HD23	1:A:89:LEU:CD2	1.95	0.96
1:A:63:ARG:NH1	1:A:63:ARG:HG2	1.73	0.96
2:B:157:ALA:O	2:B:158:ILE:HD13	1.67	0.95
2:B:142:ASP:CG	2:B:145:ARG:NH1	2.21	0.94
1:C:112:ASN:ND2	1:C:114:PHE:HE1	1.48	0.93
1:D:291:ASP:O	6:D:2133:HOH:O	1.89	0.91
1:D:267:GLU:HG3	6:D:2126:HOH:O	1.70	0.89
1:C:112:ASN:HD22	1:C:114:PHE:HE1	0.98	0.89
1:C:3:MET:CE	1:C:51:ASP:HB2	2.04	0.88
1:C:112:ASN:HD21	1:C:114:PHE:HZ	1.23	0.87
1:C:185:GLN:HE21	1:C:189:ASN:HD21	1.24	0.86
1:A:9:LEU:HD23	1:A:89:LEU:HD23	1.55	0.84
1:A:113:ILE:HD11	1:A:225:LEU:HD12	1.59	0.83
2:B:274:VAL:N	2:B:275:PRO:HD3	1.92	0.83
2:B:235:GLU:OE1	2:B:235:GLU:HA	1.79	0.82
2:B:157:ALA:C	2:B:158:ILE:HD13	1.99	0.81
1:D:91:GLN:HG3	1:D:198:LEU:HD12	1.63	0.81
2:B:141:ASN:ND2	1:C:245:GLU:O	2.14	0.80
4:D:1293:DAU:O1P	4:D:1293:DAU:O3P	1.99	0.80
4:C:1293:DAU:O1P	4:C:1293:DAU:O3P	2.00	0.79
1:A:227:THR:O	1:A:227:THR:CG2	2.31	0.78
1:D:270:ARG:NE	1:D:289:THR:HG21	1.99	0.78
1:C:34:LYS:HE2	1:C:39:TYR:CZ	2.18	0.78
2:B:275:PRO:HD2	2:B:276:LEU:H	1.50	0.77
1:C:3:MET:HE3	1:C:51:ASP:CB	2.14	0.77
1:C:61:THR:HB	1:C:62:PRO:HD3	1.66	0.77
2:B:81:LYS:HD2	6:B:2035:HOH:O	1.85	0.76
1:D:2:LYS:HD2	1:D:2:LYS:N	2.00	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:16:ARG:N	1:C:16:ARG:HD2	1.97	0.76
1:D:13:SER:OG	1:D:16:ARG:NH2	2.19	0.76
1:A:231:GLN:O	1:A:235:GLU:HG3	1.86	0.76
1:C:15:THR:HG21	1:D:278:LYS:HD2	1.68	0.76
2:B:98:GLU:N	2:B:98:GLU:CD	2.38	0.75
1:D:153:LYS:HG3	6:D:2071:HOH:O	1.86	0.74
1:C:34:LYS:HE2	1:C:39:TYR:OH	1.87	0.74
1:D:225:LEU:O	6:D:2104:HOH:O	2.06	0.73
1:A:113:ILE:CG1	1:A:225:LEU:HD12	2.19	0.73
1:D:152:ASP:HB2	6:D:2071:HOH:O	1.88	0.73
1:D:2:LYS:HA	1:D:50:ARG:HB2	1.71	0.73
1:A:9:LEU:CD2	1:A:89:LEU:HD23	2.20	0.72
2:B:122:LEU:CD2	2:B:218:MET:HE2	2.19	0.72
1:C:231:GLN:OE1	6:C:2099:HOH:O	2.06	0.71
1:D:217:MET:HB3	1:D:217:MET:CE	2.21	0.71
2:B:122:LEU:HD21	2:B:218:MET:CE	2.21	0.71
1:C:15:THR:HG21	1:D:278:LYS:CE	2.22	0.69
1:C:3:MET:CE	1:C:51:ASP:CB	2.71	0.68
1:D:160:LEU:O	1:D:205:ARG:NH1	2.13	0.68
2:B:274:VAL:N	2:B:275:PRO:CD	2.58	0.66
1:C:258:ALA:HB1	1:C:264:ILE:HD13	1.77	0.66
1:C:9:LEU:HG	4:C:1293:DAU:O21	1.96	0.66
1:A:89:LEU:HD21	1:A:109:LEU:HD21	1.78	0.65
1:C:258:ALA:CB	1:C:264:ILE:HD13	2.27	0.65
1:D:228:GLY:O	6:D:2107:HOH:O	2.13	0.65
1:C:185:GLN:HE21	1:C:189:ASN:ND2	1.93	0.65
1:D:217:MET:HG2	6:D:2097:HOH:O	1.96	0.65
1:C:63:ARG:HH11	1:C:63:ARG:HG3	1.62	0.65
2:B:138:TYR:HD1	2:B:218:MET:HG2	1.60	0.65
1:A:63:ARG:NH1	1:A:63:ARG:CG	2.35	0.64
1:A:89:LEU:HD21	1:A:109:LEU:CD2	2.27	0.64
1:D:150:GLU:HB3	1:D:159:SER:HB3	1.79	0.64
1:C:15:THR:HG21	1:D:278:LYS:CD	2.27	0.64
1:C:37:ILE:O	1:C:40:PRO:HD2	1.97	0.64
1:C:15:THR:HG21	1:D:278:LYS:HE3	1.80	0.64
1:A:113:ILE:HG12	1:A:225:LEU:HD12	1.79	0.64
1:D:217:MET:HB3	1:D:217:MET:HE3	1.79	0.63
2:B:9:LEU:HG	4:B:1293:DAU:O21	1.98	0.63
2:B:27:GLN:H	2:B:27:GLN:HE21	1.47	0.63
1:D:217:MET:CB	1:D:217:MET:CE	2.77	0.63
1:A:121:LYS:HG3	6:A:2049:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:THR:O	1:A:227:THR:HG22	1.98	0.62
2:B:275:PRO:HD2	2:B:276:LEU:N	2.15	0.62
1:A:113:ILE:HD11	1:A:225:LEU:CD1	2.29	0.62
1:A:285:LEU:O	1:A:288:MET:HG3	2.00	0.62
2:B:275:PRO:CD	2:B:276:LEU:H	2.08	0.62
6:A:2020:HOH:O	2:B:22:MET:HE1	1.99	0.61
2:B:122:LEU:HD21	2:B:218:MET:HE2	1.80	0.61
1:C:34:LYS:HE3	1:C:251:VAL:O	2.00	0.61
1:C:58:PRO:HB3	6:C:2010:HOH:O	2.01	0.61
2:B:285:LEU:O	2:B:288:GLN:HG2	2.01	0.61
1:A:27:GLN:H	1:A:27:GLN:NE2	1.99	0.60
1:A:89:LEU:CD2	1:A:109:LEU:HD21	2.32	0.60
1:A:111:ASP:HB2	1:A:227:THR:HG21	1.83	0.59
2:B:122:LEU:CD2	2:B:218:MET:CE	2.80	0.59
2:B:205:ARG:O	2:B:209:GLU:HG3	2.02	0.59
1:A:119:LEU:HB3	1:A:120:PRO:HD3	1.84	0.59
1:C:280:ASN:HB3	6:C:2107:HOH:O	2.01	0.59
1:A:98:GLU:CG	6:A:2042:HOH:O	2.50	0.58
1:A:27:GLN:H	1:A:27:GLN:HE21	1.51	0.58
1:D:205:ARG:NH1	6:D:2094:HOH:O	2.33	0.58
1:A:81:LYS:HE2	6:A:2036:HOH:O	2.03	0.58
1:A:59:GLN:HG3	6:A:2026:HOH:O	2.02	0.58
2:B:142:ASP:CG	2:B:145:ARG:HH12	1.91	0.58
1:D:266:VAL:HG23	6:D:2126:HOH:O	2.03	0.58
1:A:9:LEU:HD23	1:A:89:LEU:HD22	1.82	0.57
1:C:160:LEU:O	1:C:205:ARG:NH2	2.36	0.57
1:A:142:ASP:N	1:A:142:ASP:OD1	2.37	0.57
1:D:290:LYS:O	1:D:291:ASP:HB2	2.04	0.57
1:D:270:ARG:HD3	1:D:286:TYR:CE1	2.40	0.57
1:A:98:GLU:HG3	6:A:2042:HOH:O	2.04	0.57
1:C:238:ASN:ND2	1:D:238:ASN:HD21	2.03	0.57
1:C:15:THR:CG2	1:D:278:LYS:HE3	2.35	0.56
2:B:275:PRO:CD	2:B:276:LEU:N	2.67	0.56
1:D:2:LYS:O	1:D:51:ASP:HB2	2.05	0.56
2:B:273:ALA:C	2:B:275:PRO:CD	2.74	0.56
2:B:85:SER:HA	6:B:2033:HOH:O	2.05	0.56
2:B:138:TYR:CD1	2:B:218:MET:HG2	2.39	0.56
2:B:43:THR:OG1	6:B:2022:HOH:O	2.18	0.55
2:B:89:LEU:HD13	2:B:109:LEU:HD21	1.87	0.55
1:D:91:GLN:HG3	1:D:198:LEU:CD1	2.35	0.55
1:C:3:MET:HE3	1:C:51:ASP:CA	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:ASP:OD2	1:A:145:ARG:HD3	2.06	0.55
2:B:4:ARG:NH2	2:B:179:TYR:O	2.39	0.55
1:C:259:PHE:HE1	1:C:266:VAL:HG22	1.72	0.55
1:A:57:THR:HB	1:A:58:PRO:HD2	1.89	0.55
1:C:59:GLN:HG3	6:C:2025:HOH:O	2.08	0.54
2:B:122:LEU:HD21	2:B:218:MET:HE3	1.88	0.54
2:B:105:CYS:HB2	6:B:2039:HOH:O	2.06	0.54
1:A:227:THR:O	1:A:227:THR:HG23	2.05	0.54
1:A:81:LYS:CE	6:A:2036:HOH:O	2.55	0.54
1:A:280:ASN:H	2:B:16:ARG:NH2	2.06	0.54
2:B:57:THR:HB	2:B:58:PRO:HD2	1.89	0.54
1:C:63:ARG:HH11	1:C:63:ARG:CG	2.21	0.53
1:A:113:ILE:CD1	1:A:225:LEU:CD1	2.74	0.53
1:A:163:LYS:NZ	4:A:1293:DAU:O3P	2.38	0.53
1:A:98:GLU:HG3	1:A:99:PHE:N	2.22	0.53
1:A:113:ILE:HD13	1:A:225:LEU:HD12	1.81	0.53
1:D:266:VAL:HG23	1:D:267:GLU:N	2.24	0.53
1:C:42:SER:OG	1:C:252:SER:HB3	2.08	0.53
1:D:266:VAL:CG2	6:D:2126:HOH:O	2.57	0.53
2:B:156:THR:HB	2:B:158:ILE:HD11	1.90	0.53
1:A:4:ARG:O	1:A:49:ILE:HG23	2.08	0.52
1:C:15:THR:CG2	1:D:278:LYS:HD2	2.39	0.52
1:D:150:GLU:CD	1:D:158:ILE:HD11	2.30	0.52
1:D:89:LEU:HD13	1:D:109:LEU:HD21	1.91	0.52
1:A:284:TYR:O	1:A:288:MET:HG2	2.08	0.52
1:C:266:VAL:CG1	1:C:289:THR:HB	2.39	0.52
2:B:27:GLN:H	2:B:27:GLN:NE2	2.06	0.52
1:C:3:MET:HE1	1:C:77:ASN:HB3	1.92	0.52
2:B:7:ILE:HD11	2:B:100:ILE:HD11	1.92	0.52
2:B:135:VAL:CG2	2:B:215:VAL:HG22	2.40	0.51
1:A:142:ASP:OD2	1:A:145:ARG:CD	2.58	0.51
1:C:259:PHE:CE1	1:C:266:VAL:HG22	2.46	0.51
1:D:217:MET:CB	1:D:217:MET:HE3	2.39	0.51
1:C:160:LEU:HD11	1:C:173:VAL:HG21	1.94	0.50
2:B:142:ASP:CG	2:B:145:ARG:HH11	2.04	0.50
1:C:185:GLN:NE2	1:C:189:ASN:HD21	2.00	0.50
1:C:61:THR:N	1:C:62:PRO:CD	2.74	0.50
1:D:270:ARG:HD3	1:D:286:TYR:CZ	2.47	0.50
1:A:152:ASP:CB	1:A:156:THR:H	2.25	0.50
2:B:9:LEU:HD23	2:B:89:LEU:HD22	1.92	0.49
2:B:57:THR:HB	2:B:58:PRO:CD	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ASN:ND2	6:A:2074:HOH:O	2.44	0.48
1:C:39:TYR:N	1:C:40:PRO:CD	2.76	0.48
2:B:136:PHE:HB2	2:B:174:THR:HG22	1.94	0.48
2:B:273:ALA:C	2:B:275:PRO:HD3	2.31	0.48
1:D:59:GLN:HB3	6:D:2029:HOH:O	2.12	0.48
2:B:94:ILE:N	2:B:94:ILE:HD13	2.28	0.48
1:A:260:ARG:NH2	6:A:2126:HOH:O	2.43	0.48
2:B:3:MET:O	2:B:103:ASP:HB3	2.13	0.48
2:B:94:ILE:HD12	2:B:188:LYS:CG	2.43	0.48
1:C:115:TYR:C	1:C:115:TYR:CD2	2.86	0.48
2:B:63:ARG:HD3	6:B:2028:HOH:O	2.12	0.48
1:C:119:LEU:HB3	1:C:120:PRO:HD3	1.95	0.48
1:D:4:ARG:NE	6:D:2006:HOH:O	2.47	0.48
1:D:165:LEU:HD23	1:D:165:LEU:N	2.29	0.48
1:D:217:MET:CG	1:D:217:MET:CE	2.89	0.47
1:A:61:THR:N	1:A:62:PRO:CD	2.76	0.47
1:D:270:ARG:CZ	1:D:289:THR:HG21	2.43	0.47
1:A:239:PHE:CE2	1:A:243:ILE:HD11	2.49	0.47
1:D:112:ASN:HB3	1:D:114:PHE:CE1	2.50	0.46
1:A:4:ARG:NH1	6:A:2003:HOH:O	2.47	0.46
1:D:278:LYS:HB2	6:D:2128:HOH:O	2.16	0.46
1:D:270:ARG:HE	1:D:289:THR:HG21	1.77	0.46
1:A:57:THR:HB	1:A:58:PRO:CD	2.45	0.46
2:B:277:ILE:HD12	2:B:277:ILE:HA	1.76	0.46
1:A:161:GLU:OE1	1:A:168:LYS:HE2	2.15	0.46
1:C:258:ALA:HB3	1:C:264:ILE:CD1	2.45	0.46
1:A:278:LYS:HA	1:A:278:LYS:HD2	1.56	0.46
1:A:31:ILE:HB	1:A:39:TYR:CE1	2.51	0.46
2:B:158:ILE:HG22	2:B:158:ILE:O	2.15	0.46
1:C:63:ARG:CG	1:C:63:ARG:NH1	2.76	0.46
1:A:9:LEU:HG	4:A:1293:DAU:O21	2.16	0.46
1:A:13:SER:OG	1:A:14:GLY:N	2.49	0.46
1:C:3:MET:CE	1:C:51:ASP:CA	2.95	0.45
1:A:26:LYS:HE2	1:A:227:THR:CG2	2.46	0.45
1:D:233:LEU:HA	1:D:233:LEU:HD23	1.83	0.45
1:D:191:LYS:HD2	1:D:191:LYS:HA	1.52	0.45
2:B:136:PHE:HB2	2:B:174:THR:CG2	2.47	0.45
1:C:22:MET:HB3	1:C:22:MET:HE2	1.94	0.44
2:B:235:GLU:OE1	2:B:235:GLU:CA	2.54	0.44
1:D:2:LYS:N	1:D:2:LYS:CD	2.78	0.44
1:D:220:ARG:HD2	1:D:220:ARG:HH11	1.62	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:GLN:HG2	1:C:80:TYR:CD1	2.51	0.44
6:B:2048:HOH:O	1:C:246:ARG:HG2	2.17	0.44
2:B:61:THR:N	2:B:62:PRO:CD	2.80	0.44
1:D:57:THR:HB	1:D:58:PRO:HD2	1.99	0.44
2:B:272:LEU:O	2:B:275:PRO:CD	2.66	0.44
1:A:185:GLN:HG2	6:A:2085:HOH:O	2.17	0.44
1:D:143:PRO:HG3	1:D:170:ASN:HA	1.99	0.44
1:D:61:THR:HB	1:D:62:PRO:HD3	2.00	0.44
1:C:274:VAL:N	1:C:275:PRO:CD	2.81	0.44
1:A:152:ASP:HB3	1:A:156:THR:H	1.83	0.44
1:C:9:LEU:HD23	1:C:89:LEU:HD22	2.00	0.43
1:D:3:MET:HA	1:D:3:MET:CE	2.47	0.43
1:D:162:GLU:OE2	1:D:163:LYS:HE3	2.19	0.43
1:D:9:LEU:HD23	1:D:89:LEU:HD22	2.01	0.43
6:C:2099:HOH:O	1:D:245:GLU:HG3	2.18	0.43
2:B:85:SER:HB2	2:B:86:PRO:CD	2.48	0.43
1:A:268:GLN:HA	1:A:268:GLN:OE1	2.19	0.43
1:C:37:ILE:C	1:C:40:PRO:HD2	2.39	0.43
2:B:55:ILE:HA	2:B:81:LYS:O	2.18	0.43
1:C:258:ALA:HB3	1:C:264:ILE:HD13	2.00	0.43
1:D:28:LEU:HA	1:D:28:LEU:HD23	1.76	0.43
2:B:220:ARG:HD2	2:B:220:ARG:HH11	1.72	0.43
1:D:163:LYS:N	1:D:164:PRO:CD	2.81	0.43
1:C:238:ASN:ND2	1:D:234:ILE:CG2	2.82	0.43
2:B:74:TRP:NE1	2:B:255:GLU:HG3	2.34	0.43
1:C:16:ARG:N	1:C:16:ARG:CD	2.52	0.43
2:B:94:ILE:HD12	2:B:188:LYS:HG3	2.00	0.43
1:D:260:ARG:HG2	1:D:260:ARG:HH11	1.84	0.43
1:C:3:MET:HE3	1:C:51:ASP:N	2.34	0.42
1:D:104:ASP:OD1	1:D:180:ASP:HA	2.18	0.42
2:B:9:LEU:CD2	2:B:89:LEU:HD22	2.49	0.42
1:A:119:LEU:N	1:A:120:PRO:HD2	2.34	0.42
1:C:136:PHE:CD2	1:C:216:ALA:HB3	2.54	0.42
1:C:83:GLN:HA	1:C:84:PRO:HD2	1.81	0.42
2:B:252:SER:O	2:B:254:PRO:HD3	2.19	0.42
1:D:114:PHE:HD2	1:D:218:MET:HE1	1.84	0.42
1:D:256:GLU:HG3	1:D:288:MET:SD	2.59	0.42
2:B:7:ILE:HD12	2:B:105:CYS:SG	2.60	0.41
1:C:15:THR:HB	1:C:16:ARG:HH11	1.84	0.41
1:C:57:THR:HB	1:C:58:PRO:HD2	2.02	0.41
1:C:2:LYS:HD2	1:C:2:LYS:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:180:ASP:OD1	2:B:182:ASP:HB2	2.20	0.41
1:C:114:PHE:CD2	1:C:218:MET:HE3	2.55	0.41
2:B:100:ILE:HG22	2:B:100:ILE:O	2.18	0.41
2:B:94:ILE:HD12	2:B:188:LYS:HG2	2.03	0.41
1:A:152:ASP:HB2	1:A:156:THR:H	1.86	0.41
2:B:100:ILE:CG2	2:B:100:ILE:O	2.68	0.41
1:D:270:ARG:NH2	1:D:289:THR:HG22	2.35	0.41
1:A:72:SER:HB3	6:A:2032:HOH:O	2.20	0.41
1:D:81:LYS:HE2	5:D:1294:SO4:O2	2.20	0.41
1:D:16:ARG:CD	1:D:16:ARG:H	2.33	0.41
2:B:73:GLN:HG2	2:B:74:TRP:CD1	2.57	0.41
1:A:119:LEU:N	1:A:120:PRO:CD	2.84	0.40
1:A:89:LEU:CD2	1:A:109:LEU:CD2	2.94	0.40
2:B:126:ALA:HB2	2:B:134:THR:HG21	2.03	0.40
1:C:234:ILE:HG21	1:C:234:ILE:HD13	1.86	0.40
1:A:116:GLY:HA3	1:A:222:TYR:CD2	2.56	0.40
1:A:200:ILE:HA	1:A:200:ILE:HD12	1.96	0.40
1:C:126:ALA:HB2	1:C:134:THR:HG21	2.03	0.40

All (2) 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)
6:A:2056:HOH:O	6:C:2051:HOH:O[1_554]	2.12	0.08
6:B:2002:HOH:O	6:D:2133:HOH:O[2_747]	2.16	0.04

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	288/293 (98%)	284 (99%)	3 (1%)	1 (0%)	46 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	288/293 (98%)	284 (99%)	4 (1%)	0	100	100
1	D	288/293 (98%)	284 (99%)	2 (1%)	2 (1%)	26	14
2	B	288/293 (98%)	282 (98%)	6 (2%)	0	100	100
All	All	1152/1172 (98%)	1134 (98%)	15 (1%)	3 (0%)	46	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	D	32	TYR
1	D	124	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	244/247 (99%)	228 (93%)	16 (7%)	21	10
1	C	242/247 (98%)	229 (95%)	13 (5%)	27	15
1	D	243/247 (98%)	236 (97%)	7 (3%)	50	40
2	B	242/247 (98%)	230 (95%)	12 (5%)	30	18
All	All	971/988 (98%)	923 (95%)	48 (5%)	31	18

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	16	ARG
1	A	27	GLN
1	A	63	ARG
1	A	79	GLN
1	A	81	LYS
1	A	98	GLU
1	A	112	ASN

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Mol	Chain	Res	Type
1	A	113	ILE
1	A	158	ILE
1	A	160	LEU
1	A	227	THR
1	A	246	ARG
1	A	280	ASN
1	A	288	MET
1	A	290	LYS
2	B	4	ARG
2	B	27	GLN
2	B	41	LEU
2	B	98	GLU
2	B	112	ASN
2	B	130	GLU
2	B	153	LYS
2	B	208	LEU
2	B	220	ARG
2	B	246	ARG
2	B	267	GLU
2	B	280	ASN
1	C	2	LYS
1	C	16	ARG
1	C	22	MET
1	C	24	VAL
1	C	27	GLN
1	C	79	GLN
1	C	115	TYR
1	C	190	LEU
1	C	195	ARG
1	C	264	ILE
1	C	268	GLN
1	C	277	ILE
1	C	290	LYS
1	D	3	MET
1	D	16	ARG
1	D	41	LEU
1	D	59	GLN
1	D	79	GLN
1	D	118	ASP
1	D	119	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	128	ASN
1	A	170	ASN
1	A	231	GLN
1	A	280	ASN
1	A	283	GLN
2	B	27	GLN
2	B	154	ASN
1	C	27	GLN
1	C	79	GLN
1	C	189	ASN
1	C	230	HIS
1	C	238	ASN
1	C	268	GLN
1	D	27	GLN
1	D	59	GLN

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 ⓘ

9 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
3	TYD	A	1292	-	19,26,26	1.08	2 (10%)	27,40,40	2.25	5 (18%)
4	DAU	A	1293	-	19,28,38	1.52	3 (15%)	29,40,58	3.05	8 (27%)
3	TYD	B	1292	-	19,26,26	1.08	1 (5%)	27,40,40	1.72	3 (11%)
4	DAU	B	1293	-	19,28,38	1.57	3 (15%)	29,40,58	3.18	12 (41%)
3	TYD	C	1292	-	19,26,26	1.15	1 (5%)	27,40,40	2.08	3 (11%)
4	DAU	C	1293	-	19,28,38	1.52	3 (15%)	29,40,58	3.16	8 (27%)
3	TYD	D	1292	-	19,26,26	1.31	2 (10%)	27,40,40	2.17	6 (22%)
4	DAU	D	1293	1	19,28,38	1.47	3 (15%)	29,40,58	3.02	9 (31%)
5	SO4	D	1294	-	4,4,4	0.17	0	6,6,6	0.30	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
3	TYD	A	1292	-	-	0/12/28/28	0/2/2/2
4	DAU	A	1293	-	-	0/10/32/55	0/2/2/3
3	TYD	B	1292	-	-	0/12/28/28	0/2/2/2
4	DAU	B	1293	-	-	0/10/32/55	0/2/2/3
3	TYD	C	1292	-	-	0/12/28/28	0/2/2/2
4	DAU	C	1293	-	-	0/10/32/55	0/2/2/3
3	TYD	D	1292	-	-	0/12/28/28	0/2/2/2
4	DAU	D	1293	1	-	0/10/32/55	0/2/2/3
5	SO4	D	1294	-	-	0/0/0/0	0/0/0/0

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1292	TYD	C6-N1	-2.17	1.31	1.35
3	A	1292	TYD	PA-O2A	-2.16	1.45	1.54
4	A	1293	DAU	O41-C41	2.01	1.29	1.24
4	C	1293	DAU	O41-C41	2.02	1.29	1.24
4	D	1293	DAU	O41-C41	2.04	1.29	1.24
4	B	1293	DAU	O41-C41	2.05	1.29	1.24
4	C	1293	DAU	P2-O4P	2.48	1.65	1.54
4	D	1293	DAU	P2-O4P	2.95	1.67	1.54
4	A	1293	DAU	P2-O4P	3.20	1.68	1.54
4	B	1293	DAU	P2-O4P	3.24	1.68	1.54
3	B	1292	TYD	O4-C4	3.31	1.32	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1292	TYD	O4-C4	3.43	1.32	1.24
3	C	1292	TYD	O4-C4	3.65	1.33	1.24
3	D	1292	TYD	O4-C4	3.99	1.34	1.24
4	A	1293	DAU	P2-O3P	4.46	1.67	1.51
4	D	1293	DAU	P2-O3P	4.66	1.68	1.51
4	C	1293	DAU	P2-O3P	5.03	1.69	1.51
4	B	1293	DAU	P2-O3P	5.22	1.70	1.51

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1293	DAU	N11-C21-N31	-10.18	121.84	128.33
4	A	1293	DAU	N11-C21-N31	-10.14	121.86	128.33
4	B	1293	DAU	N11-C21-N31	-10.13	121.87	128.33
4	C	1293	DAU	N11-C21-N31	-10.12	121.88	128.33
4	A	1293	DAU	C51-C61-N11	-9.29	116.94	125.23
4	C	1293	DAU	C51-C61-N11	-9.28	116.96	125.23
4	B	1293	DAU	C51-C61-N11	-9.21	117.02	125.23
4	D	1293	DAU	C51-C61-N11	-9.20	117.02	125.23
4	C	1293	DAU	O5-C1-O1	-6.41	102.91	111.36
3	D	1292	TYD	C5-C4-N3	-5.62	118.88	125.14
3	A	1292	TYD	C5-C4-N3	-5.47	119.05	125.14
3	C	1292	TYD	C5-C4-N3	-5.05	119.52	125.14
3	B	1292	TYD	C5-C4-N3	-4.55	120.07	125.14
3	D	1292	TYD	PA-O3A-PB	-3.45	121.10	132.67
4	A	1293	DAU	C2-C3-C4	-3.42	104.40	110.79
3	A	1292	TYD	PA-O3A-PB	-3.41	121.24	132.67
4	B	1293	DAU	C2-C3-C4	-3.26	104.71	110.79
4	D	1293	DAU	O2-C2-C1	-2.95	103.55	110.02
4	B	1293	DAU	O3-C3-C4	-2.87	103.88	110.34
4	D	1293	DAU	C2-C3-C4	-2.62	105.90	110.79
4	A	1293	DAU	C51-C41-N31	-2.53	122.32	125.14
4	C	1293	DAU	C51-C41-N31	-2.52	122.34	125.14
4	B	1293	DAU	C51-C41-N31	-2.48	122.38	125.14
4	D	1293	DAU	C51-C41-N31	-2.42	122.44	125.14
3	D	1292	TYD	O3B-PB-O1B	-2.33	103.08	110.58
4	B	1293	DAU	O2-C2-C1	-2.23	105.13	110.02
4	D	1293	DAU	O3-C3-C4	-2.09	105.63	110.34
4	A	1293	DAU	C6-C5-C4	-2.07	107.92	113.02
3	D	1292	TYD	C5M-C5-C4	2.02	122.66	120.05
4	C	1293	DAU	O1-P2-O3P	2.09	117.69	109.46
3	C	1292	TYD	C5M-C5-C6	2.14	122.92	118.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1293	DAU	O3-C3-C2	2.20	115.28	110.34
4	B	1293	DAU	O5-C5-C4	2.20	113.81	109.68
4	B	1293	DAU	C3-C4-C5	2.46	114.49	110.20
3	B	1292	TYD	O2B-PB-O3A	2.48	116.35	105.09
4	C	1293	DAU	O5-C5-C4	2.49	114.35	109.68
3	A	1292	TYD	O3A-PA-O5'	2.58	109.77	102.94
4	D	1293	DAU	OPP-P2-O1	2.62	111.26	103.58
4	B	1293	DAU	O1-C1-C2	2.64	113.33	108.39
4	B	1293	DAU	C41-N31-C21	2.70	117.58	115.25
4	D	1293	DAU	C41-N31-C21	2.72	117.60	115.25
3	A	1292	TYD	O2B-PB-O3A	2.74	117.51	105.09
4	C	1293	DAU	C41-N31-C21	2.75	117.63	115.25
4	A	1293	DAU	C41-N31-C21	2.78	117.65	115.25
4	B	1293	DAU	C61-N11-C21	3.75	121.50	115.47
4	C	1293	DAU	C61-N11-C21	3.76	121.52	115.47
4	D	1293	DAU	C61-N11-C21	3.77	121.53	115.47
4	A	1293	DAU	C61-N11-C21	3.78	121.55	115.47
4	B	1293	DAU	C1-O5-C5	3.86	121.24	113.75
3	D	1292	TYD	O3A-PA-O5'	5.02	116.25	102.94
3	B	1292	TYD	C4-N3-C2	5.05	119.61	115.25
3	D	1292	TYD	C4-N3-C2	5.92	120.36	115.25
3	A	1292	TYD	C4-N3-C2	8.01	122.17	115.25
3	C	1292	TYD	C4-N3-C2	8.23	122.36	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1293	DAU	2	0
4	B	1293	DAU	1	0
4	C	1293	DAU	2	0
4	D	1293	DAU	1	0
5	D	1294	SO4	1	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 ⓘ

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.