



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:59 AM UTC

PDB ID : 3N98 / pdb_00003n98
Title : Crystal structure of TK1436, a GH57 branching enzyme from hyperthermophilic archaeon *Thermococcus kodakaraensis*, in complex with glucose and additives
Authors : Santos, C.R.; Tonoli, C.C.C.; Trindade, D.M.; Betzel, C.; Takata, H.; Kuriki, T.; Kanai, T.; Imanaka, T.; Arni, R.K.; Murakami, M.T.
Deposited on : 2010-05-28
Resolution : 1.87 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

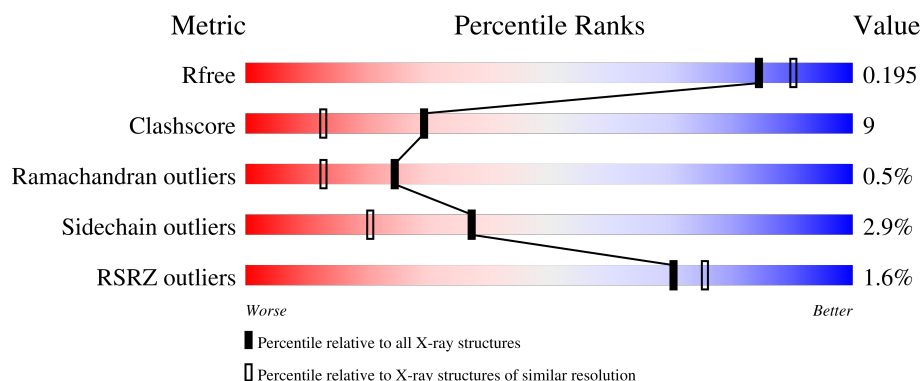
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	562	

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
2	BGC	A	564	-	-	X	-

2 Entry composition [i](#)

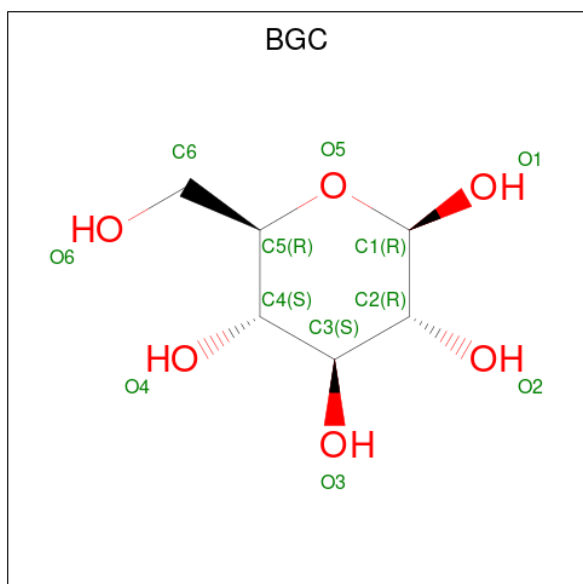
There are 7 unique types of molecules in this entry. The entry contains 5128 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 alpha-amylase, GH57 family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	550	4694	3033	796	853	12	0	25	0

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



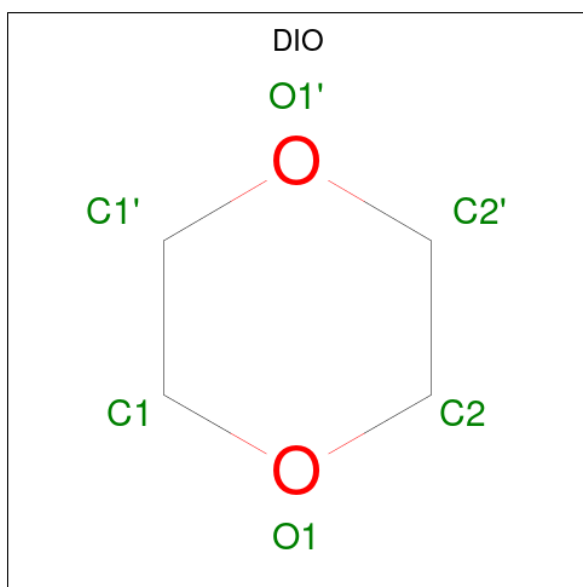
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 1,4-DIETHYLENE DIOXIDE (CCD ID: DIO) (formula: C₄H₈O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	4	2		

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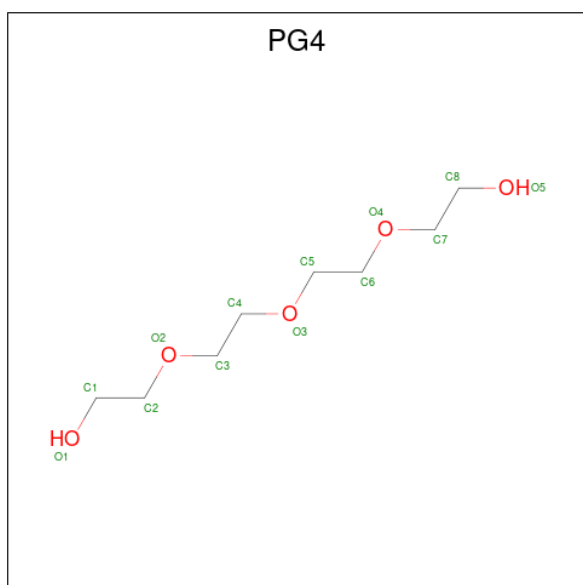
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	4	2		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			13	8	5		

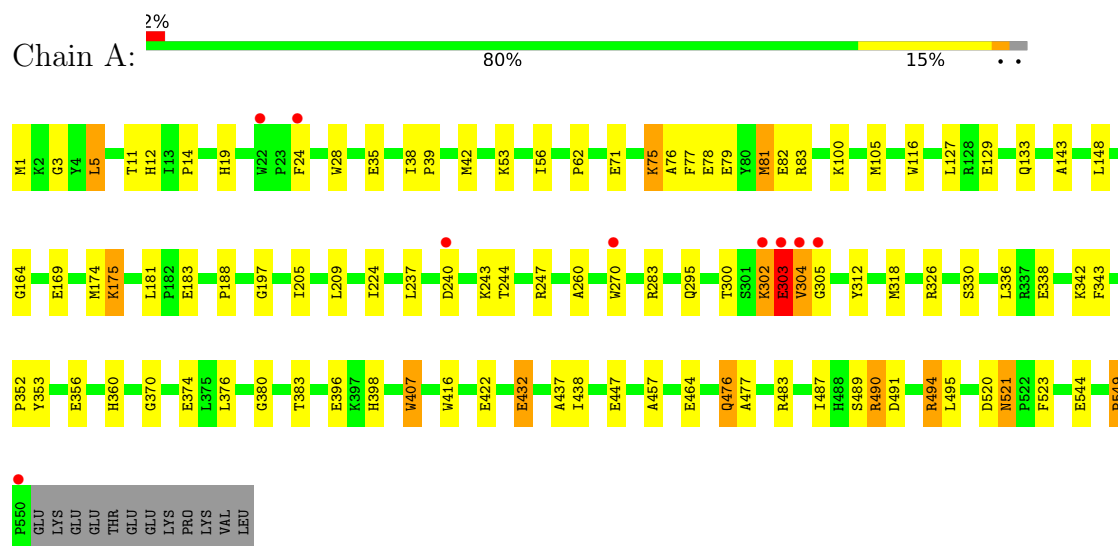
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	348	Total	O	0	0
			348	348		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: alpha-amylase, GH57 family



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.67Å 79.02Å 134.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.08 – 1.87 68.08 – 1.87	Depositor EDS
% Data completeness (in resolution range)	95.4 (68.08-1.87) 95.4 (68.08-1.87)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.159 , 0.196 0.157 , 0.195	Depositor DCC
R_{free} test set	3024 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 58.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5128	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, BGC, PEG, DIO, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.61	20/4945 (0.4%)	1.30	22/6689 (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	A	0	1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	143	ALA	CA-CB	6.38	1.63	1.53
1	A	237	LEU	N-CA	6.34	1.53	1.45
1	A	303	GLU	N-CA	6.22	1.54	1.46
1	A	76	ALA	CA-C	6.21	1.60	1.52
1	A	304	VAL	CA-CB	6.17	1.62	1.54
1	A	549	PRO	CA-C	5.99	1.55	1.51
1	A	437	ALA	N-CA	5.80	1.53	1.46
1	A	12	HIS	C-O	5.75	1.30	1.23
1	A	244	THR	CA-C	5.73	1.59	1.52
1	A	457	ALA	N-CA	5.70	1.53	1.46
1	A	56	ILE	N-CA	5.70	1.53	1.46
1	A	300	THR	CA-CB	5.57	1.62	1.53
1	A	330	SER	N-CA	5.44	1.52	1.46
1	A	336	LEU	CA-C	5.44	1.59	1.52
1	A	3	GLY	N-CA	5.42	1.51	1.44
1	A	477	ALA	CA-CB	5.35	1.57	1.53
1	A	520	ASP	N-CA	5.30	1.53	1.46
1	A	76	ALA	C-O	5.24	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	PRO	N-CA	-5.15	1.41	1.47
1	A	56	ILE	CA-CB	-5.13	1.50	1.55

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	544	GLU	CA-C-N	-8.51	111.61	120.38
1	A	544	GLU	C-N-CA	-8.51	111.61	120.38
1	A	181	LEU	N-CA-C	-7.62	100.49	109.93
1	A	521	ASN	CA-C-N	-6.45	113.33	119.85
1	A	521	ASN	C-N-CA	-6.45	113.33	119.85
1	A	71	GLU	N-CA-C	6.43	118.83	111.11
1	A	24	PHE	CA-C-N	-6.12	115.55	121.31
1	A	24	PHE	C-N-CA	-6.12	115.55	121.31
1	A	477	ALA	N-CA-C	-6.06	97.79	109.24
1	A	432	GLU	N-CA-C	-6.04	104.70	111.28
1	A	380	GLY	CA-C-N	-6.00	115.37	122.93
1	A	380	GLY	C-N-CA	-6.00	115.37	122.93
1	A	148	LEU	CA-C-N	-5.66	113.78	119.56
1	A	148	LEU	C-N-CA	-5.66	113.78	119.56
1	A	490	ARG	NE-CZ-NH1	5.48	126.98	121.50
1	A	82	GLU	CA-CB-CG	-5.47	103.15	114.10
1	A	476	GLN	CA-C-N	-5.22	112.33	122.62
1	A	476	GLN	C-N-CA	-5.22	112.33	122.62
1	A	376	LEU	N-CA-C	-5.22	105.67	111.36
1	A	489	SER	N-CA-C	-5.20	105.69	111.36
1	A	438	ILE	N-CA-C	5.07	115.30	110.53
1	A	197	GLY	N-CA-C	5.03	120.71	114.37

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	305	GLY	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4694	0	4551	75	0
2	A	36	0	36	11	0
3	A	18	0	24	2	0
4	A	12	0	16	3	0
5	A	7	0	10	0	0
6	A	13	0	18	3	0
7	A	348	0	0	26	0
All	All	5128	0	4655	82	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 9.

All (82) 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:270[B]:TRP:HZ2	7:A:911:HOH:O	1.21	1.20
1:A:169[B]:GLU:OE2	1:A:175:LYS:HG2	1.46	1.13
1:A:270[B]:TRP:HZ3	7:A:617:HOH:O	1.34	1.08
1:A:483[A]:ARG:NH1	1:A:549:PRO:O	1.90	1.03
1:A:270[B]:TRP:CZ2	7:A:911:HOH:O	2.02	0.99
6:A:572:PG4:H61	7:A:880:HOH:O	1.63	0.97
1:A:247[B]:ARG:NH1	1:A:398:HIS:CE1	2.33	0.95
1:A:169[B]:GLU:OE2	1:A:175:LYS:CG	2.16	0.93
1:A:270[B]:TRP:CZ3	7:A:617:HOH:O	2.14	0.92
1:A:247[B]:ARG:HH11	1:A:398:HIS:CE1	1.86	0.91
3:A:567:GOL:O3	7:A:911:HOH:O	1.79	0.89
1:A:476:GLN:HE22	2:A:564:BGC:C6	1.84	0.89
1:A:129[B]:GLU:HG2	1:A:133:GLN:NE2	1.95	0.80
1:A:39:PRO:HA	1:A:42[A]:MET:HE2	1.68	0.75
1:A:476:GLN:HE22	2:A:564:BGC:H6C1	1.51	0.75
1:A:42[A]:MET:CE	7:A:862:HOH:O	2.36	0.74
1:A:169[B]:GLU:OE2	1:A:175:LYS:CD	2.35	0.74
1:A:224:ILE:HD12	1:A:260:ALA:HB2	1.70	0.74
2:A:565:BGC:O6	7:A:868:HOH:O	2.05	0.73
1:A:39:PRO:HA	1:A:42[A]:MET:CE	2.19	0.73
1:A:283[B]:ARG:HD2	1:A:295:GLN:HB3	1.71	0.72
6:A:572:PG4:H42	7:A:880:HOH:O	1.91	0.70
1:A:42[A]:MET:HE3	7:A:862:HOH:O	1.91	0.69
1:A:243[B]:LYS:HE2	7:A:913:HOH:O	1.92	0.69
1:A:77:PHE:CZ	1:A:81:MET:HG2	2.29	0.67
6:A:572:PG4:H21	7:A:693:HOH:O	1.93	0.67
1:A:53:LYS:NZ	7:A:639:HOH:O	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:563:BGC:H3	7:A:893:HOH:O	1.97	0.63
1:A:476:GLN:NE2	2:A:564:BGC:C6	2.60	0.62
1:A:79[A]:GLU:HG2	1:A:83:ARG:NH2	2.15	0.61
1:A:129[B]:GLU:CG	1:A:133:GLN:NE2	2.64	0.61
1:A:205:ILE:O	1:A:209[A]:LEU:HD13	2.01	0.60
1:A:28:TRP:CE3	2:A:564:BGC:O3	2.56	0.59
1:A:247[A]:ARG:NH2	7:A:890:HOH:O	2.35	0.59
1:A:100[A]:LYS:HD2	7:A:580:HOH:O	2.04	0.57
1:A:38[A]:ILE:HD11	1:A:116:TRP:HE3	1.71	0.56
1:A:270[B]:TRP:HH2	1:A:352:PRO:O	1.89	0.55
1:A:42[A]:MET:HE1	7:A:862:HOH:O	2.05	0.54
1:A:62:PRO:HG2	1:A:164:GLY:HA2	1.88	0.54
1:A:35[B]:GLU:OE2	1:A:360:HIS:HD2	1.91	0.54
1:A:476:GLN:NE2	2:A:564:BGC:H6C1	2.19	0.54
1:A:169[B]:GLU:OE2	1:A:175:LYS:CE	2.56	0.53
1:A:338:GLU:OE1	7:A:729:HOH:O	2.19	0.53
1:A:169[B]:GLU:OE2	1:A:175:LYS:NZ	2.42	0.53
1:A:318:MET:HG2	4:A:570:DIO:H22	1.91	0.52
1:A:129[B]:GLU:HG2	1:A:133:GLN:HE22	1.74	0.52
1:A:39:PRO:CA	1:A:42[A]:MET:HE2	2.39	0.51
1:A:283[B]:ARG:HD2	1:A:295:GLN:CB	2.39	0.51
1:A:370:GLY:O	1:A:374[A]:GLU:HG3	2.11	0.50
1:A:490:ARG:HD3	1:A:494[C]:ARG:HD2	1.94	0.49
1:A:75:LYS:HE3	1:A:78:GLU:OE1	2.13	0.49
1:A:356[B]:GLU:OE1	7:A:865:HOH:O	2.20	0.49
1:A:11:THR:HG21	4:A:569:DIO:H2'2	1.96	0.48
2:A:565:BGC:H5	7:A:918:HOH:O	2.13	0.48
1:A:247[B]:ARG:NH2	1:A:343:PHE:CD1	2.82	0.47
1:A:270[B]:TRP:CH2	1:A:352:PRO:O	2.67	0.47
1:A:432:GLU:OE2	1:A:464:GLU:OE2	2.32	0.47
1:A:476:GLN:NE2	2:A:564:BGC:H6C2	2.27	0.47
1:A:39:PRO:HA	1:A:42[A]:MET:HE3	1.95	0.47
1:A:483[A]:ARG:HG2	1:A:487:ILE:CD1	2.45	0.46
1:A:326:ARG:NE	7:A:916:HOH:O	2.47	0.46
1:A:14:PRO:HD2	2:A:564:BGC:H2	1.99	0.45
1:A:312:TYR:OH	4:A:570:DIO:H2'2	2.15	0.45
1:A:483[B]:ARG:HG3	7:A:876:HOH:O	2.15	0.45
1:A:247[B]:ARG:HH12	1:A:398:HIS:CE1	2.28	0.44
1:A:270[B]:TRP:CH2	1:A:353:TYR:CD2	3.06	0.43
1:A:476:GLN:HE22	2:A:564:BGC:H6C2	1.71	0.43
1:A:302:LYS:C	1:A:303:GLU:HG3	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:490:ARG:HD2	1:A:491:ASP:OD1	2.18	0.43
1:A:247[B]:ARG:NH1	1:A:398:HIS:HE1	2.10	0.42
1:A:129[B]:GLU:CG	1:A:133:GLN:HE22	2.31	0.42
1:A:342:LYS:HG2	1:A:343:PHE:CE2	2.54	0.42
1:A:407:TRP:CE2	3:A:567:GOL:H12	2.55	0.41
1:A:174:MET:CE	7:A:902:HOH:O	2.67	0.41
1:A:19:HIS:C	7:A:919:HOH:O	2.64	0.41
1:A:100[B]:LYS:HE3	1:A:100[B]:LYS:HB2	1.13	0.41
1:A:105[B]:MET:HE3	7:A:579:HOH:O	2.20	0.41
1:A:5:LEU:O	1:A:383:THR:HA	2.22	0.40
1:A:270[B]:TRP:CZ2	1:A:353:TYR:HD2	2.39	0.40
1:A:447:GLU:HG3	7:A:751:HOH:O	2.22	0.40
1:A:183:GLU:HG3	1:A:416:TRP:CZ2	2.57	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	573/562 (102%)	560 (98%)	10 (2%)	3 (0%)	24 13

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	304	VAL
1	A	523	PHE
1	A	521	ASN

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	502/489 (103%)	485 (97%)	17 (3%)	32	16

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	LEU
1	A	75	LYS
1	A	81	MET
1	A	127	LEU
1	A	175	LYS
1	A	240[A]	ASP
1	A	240[B]	ASP
1	A	302	LYS
1	A	303	GLU
1	A	396	GLU
1	A	407	TRP
1	A	422[A]	GLU
1	A	422[B]	GLU
1	A	494[B]	ARG
1	A	494[C]	ARG
1	A	495	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	GLN
1	A	273	HIS
1	A	360	HIS
1	A	398	HIS
1	A	476	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

10 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	$\# Z > 2$
6	PG4	A	572	-	12,12,12	0.56	0	11,11,11	0.58	0
2	BGC	A	565	-	12,12,12	0.57	0	17,17,17	1.11	1 (5%)
3	GOL	A	566	-	5,5,5	0.37	0	5,5,5	1.00	0
5	PEG	A	571	-	6,6,6	0.29	0	5,5,5	0.88	0
4	DIO	A	569	-	6,6,6	0.66	0	6,6,6	1.35	0
4	DIO	A	570	-	6,6,6	0.82	0	6,6,6	1.98	2 (33%)
3	GOL	A	567	-	5,5,5	0.49	0	5,5,5	1.56	1 (20%)
2	BGC	A	564	-	12,12,12	0.69	0	17,17,17	1.08	2 (11%)
3	GOL	A	568	-	5,5,5	1.01	0	5,5,5	0.90	0
2	BGC	A	563	-	12,12,12	0.86	0	17,17,17	1.67	6 (35%)

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
6	PG4	A	572	-	-	5/10/10/10	-
2	BGC	A	565	-	-	1/2/22/22	0/1/1/1
3	GOL	A	566	-	-	2/4/4/4	-
5	PEG	A	571	-	-	0/4/4/4	-
4	DIO	A	569	-	-	-	0/1/1/1
4	DIO	A	570	-	-	-	0/1/1/1
3	GOL	A	567	-	-	0/4/4/4	-
2	BGC	A	564	-	-	0/2/22/22	0/1/1/1
3	GOL	A	568	-	-	0/4/4/4	-
2	BGC	A	563	-	-	2/2/22/22	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	563	BGC	C1-O5-C5	3.65	120.72	113.65
4	A	570	DIO	C2-O1-C1	3.41	120.89	109.88
4	A	570	DIO	C2'-O1'-C1'	3.05	119.73	109.88
2	A	563	BGC	O5-C1-C2	2.99	115.56	110.30
2	A	563	BGC	O4-C4-C3	-2.58	104.30	110.38
2	A	563	BGC	O5-C5-C4	2.53	114.26	109.70
2	A	565	BGC	C3-C4-C5	-2.27	106.11	110.23
2	A	564	BGC	O5-C1-C2	2.20	114.17	110.30
2	A	564	BGC	O3-C3-C2	-2.10	105.44	110.38
3	A	567	GOL	O2-C2-C3	-2.10	100.50	109.18
2	A	563	BGC	O5-C5-C6	2.06	111.55	106.44
2	A	563	BGC	C3-C4-C5	2.05	113.95	110.23

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	563	BGC	O5-C5-C6-O6
2	A	563	BGC	C4-C5-C6-O6
6	A	572	PG4	O2-C3-C4-O3
3	A	566	GOL	O2-C2-C3-O3
6	A	572	PG4	O4-C7-C8-O5
3	A	566	GOL	C1-C2-C3-O3
6	A	572	PG4	C8-C7-O4-C6
2	A	565	BGC	O5-C5-C6-O6
6	A	572	PG4	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
6	A	572	PG4	O3-C5-C6-O4

There are no ring outliers.

7 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	572	PG4	3	0
2	A	565	BGC	2	0
4	A	569	DIO	1	0
4	A	570	DIO	2	0
3	A	567	GOL	2	0
2	A	564	BGC	8	0
2	A	563	BGC	1	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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	550/562 (97%)	-0.25	9 (1%) 70 75	11, 21, 37, 63	48 (8%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	304	VAL	7.6
1	A	305	GLY	4.4
1	A	22	TRP	3.1
1	A	550	PRO	2.7
1	A	302	LYS	2.7
1	A	240[A]	ASP	2.5
1	A	270[A]	TRP	2.4
1	A	24	PHE	2.2
1	A	303	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	A	565	12/12	0.70	0.19	82,85,86,88	0
2	BGC	A	564	12/12	0.79	0.16	79,82,83,84	0
4	DIO	A	570	6/6	0.82	0.16	41,43,45,50	0
3	GOL	A	566	6/6	0.86	0.12	44,49,50,52	0
2	BGC	A	563	12/12	0.86	0.14	65,70,76,77	0
6	PG4	A	572	13/13	0.88	0.14	45,49,54,59	0
3	GOL	A	568	6/6	0.89	0.10	32,37,41,42	0
3	GOL	A	567	6/6	0.91	0.10	25,46,50,55	0
4	DIO	A	569	6/6	0.91	0.12	31,34,39,43	0
5	PEG	A	571	7/7	0.95	0.09	40,42,46,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.