



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:01 AM UTC

PDB ID : 1T8E / pdb_00001t8e
Title : T7 DNA Polymerase Ternary Complex with dCTP at the Insertion Site.
Authors : Brieba, L.G.; Eichman, B.F.; Kokoska, R.J.; Doublié, S.; Kunkel, T.A.; Ellenberger, T.
Deposited on : 2004-05-12
Resolution : 2.54 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

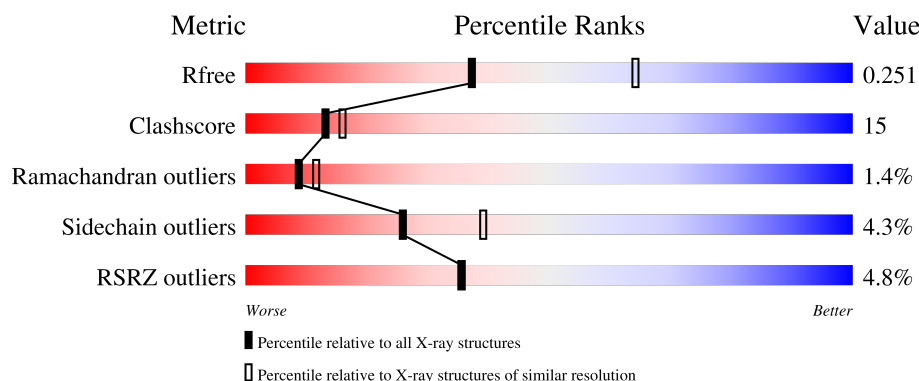
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	C	22	23% 32% 59% 9%
2	D	26	35% 54% 27% 12% • •
3	A	698	3% 71% 26% •
4	B	108	6% 55% 37% 6% •

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 7741 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 DNA chain called 5'-D(P*CP*GP*AP*AP*AP*AP*CP*GP*AP*CP*GP*GP*CP*CP*AP*GP*TP*GP*CP*CP*AP*(2DT))-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	22	Total	C	N	O	P	0	0	0
			451	213	90	126	22			

- Molecule 2 is a DNA chain called 25-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	25	Total	C	N	O	P	0	0	0
			514	244	89	156	25			

- Molecule 3 is a protein called DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	698	Total	C	N	O	S	0	0	0
			5537	3522	963	1028	24			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LYS	deletion	UNP P00581
A	?	-	ARG	deletion	UNP P00581
A	?	-	PHE	deletion	UNP P00581
A	?	-	GLY	deletion	UNP P00581
A	?	-	SER	deletion	UNP P00581
A	?	-	HIS	deletion	UNP P00581

- Molecule 4 is a protein called thioredoxin 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	106	Total	C	N	O	S	0	0	0
			798	514	128	153	3			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

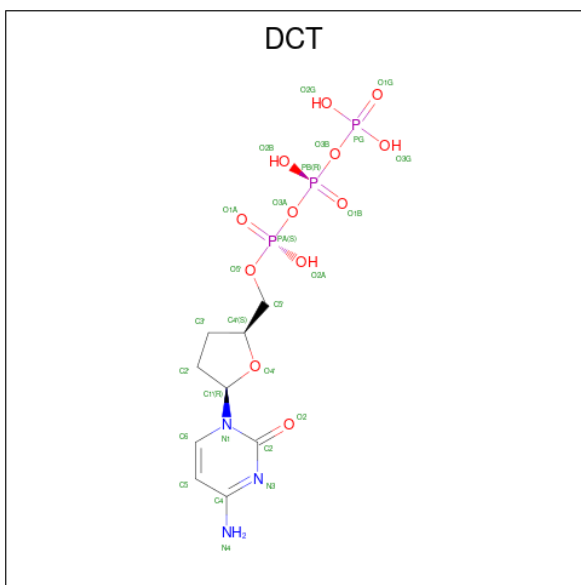
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Mg	0	0
			3	3		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



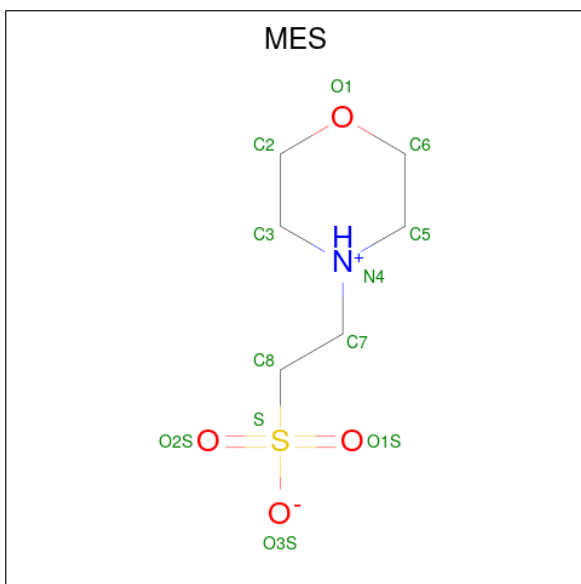
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is 2',3'-DIDEOXYCYTIDINE 5'-TRIPHOSPHATE (CCD ID: DCT) (formula: C₉H₁₆N₃O₁₂P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			27	9	3	12	3		

- Molecule 8 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $\text{C}_8\text{H}_{18}\text{O}_5$).



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

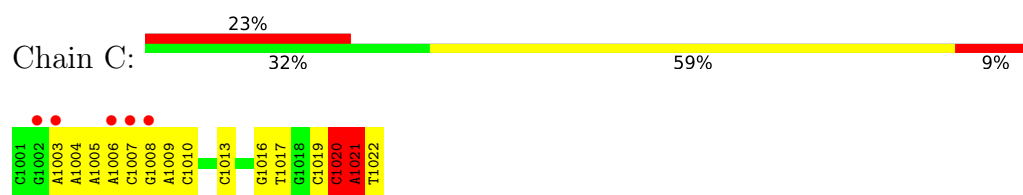
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	16	Total	O	0	0
			16	16		
10	D	27	Total	O	0	0
			27	27		
10	A	319	Total	O	0	0
			319	319		
10	B	19	Total	O	0	0
			19	19		

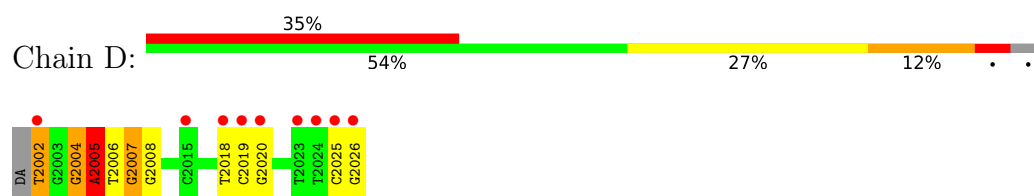
3 Residue-property plots

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.

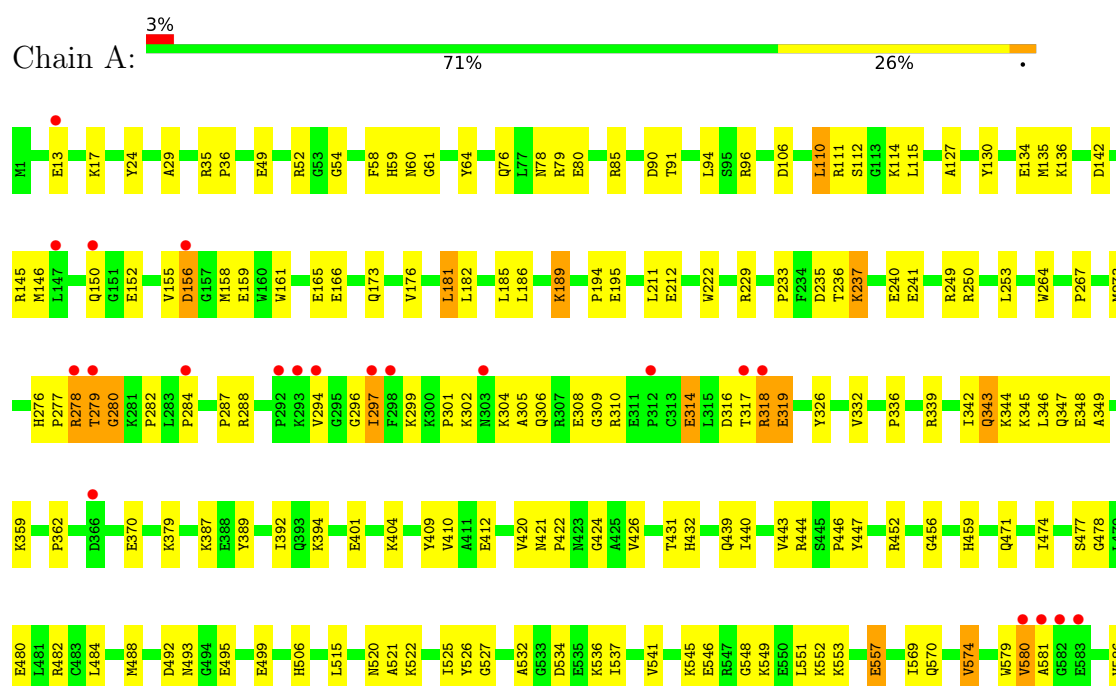
- Molecule 1: 5'-D(P*CP*GP*AP*AP*AP*AP*CP*GP*AP*CP*GP*GP*CP*CP*AP*GP*TP*GP*CP*CP*AP*(2DT))-3'

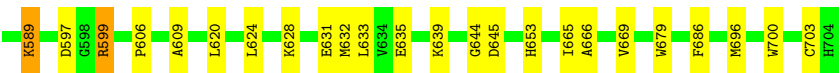


- Molecule 2: 25-MER

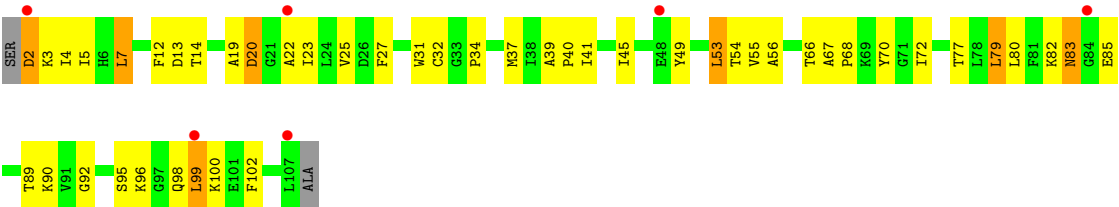


- Molecule 3: DNA polymerase





• Molecule 4: thioredoxin 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	106.28Å 218.12Å 52.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.71 – 2.54 28.71 – 2.54	Depositor EDS
% Data completeness (in resolution range)	95.0 (28.71-2.54) 94.9 (28.71-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 2.54Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.255 0.216 , 0.251	Depositor DCC
R_{free} test set	1923 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.2	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	0.92	EDS
Total number of atoms	7741	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: 2DT, DCT, MES, SO4, MG, PG4

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	0.32	0/486	0.72	2/747 (0.3%)
2	D	0.34	0/574	0.90	4/885 (0.5%)
3	A	0.41	0/5674	0.89	10/7683 (0.1%)
4	B	0.41	0/813	0.84	1/1106 (0.1%)
All	All	0.40	0/7547	0.88	17/10421 (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	C	0	2
2	D	0	3
All	All	0	5

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2002	DT	N1-C1'-C2'	7.25	124.38	113.50
3	A	127	ALA	N-CA-C	7.23	118.85	110.97
3	A	527	GLY	N-CA-C	-6.99	104.35	112.73
3	A	589	LYS	N-CA-C	-6.90	102.85	112.45
3	A	703	CYS	N-CA-C	6.42	120.72	113.02
3	A	304	LYS	N-CA-C	-6.27	104.33	112.23
3	A	586	VAL	N-CA-C	6.22	117.55	108.53
3	A	401	GLU	N-CA-C	6.11	120.94	112.45
1	C	1020	DC	N1-C1'-C2'	-5.65	105.03	113.50
2	D	2004	DG	N9-C1'-C2'	-5.59	105.11	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2007	DG	N9-C1'-C2'	-5.46	105.31	113.50
2	D	2005	DA	N9-C1'-C2'	-5.40	105.40	113.50
4	B	7	LEU	N-CA-C	5.33	116.96	110.41
3	A	574	VAL	N-CA-C	5.24	116.84	108.87
3	A	166	GLU	N-CA-C	-5.19	105.30	111.69
3	A	424	GLY	N-CA-C	5.18	120.37	114.67
1	C	1021	DA	N9-C1'-C2'	-5.18	105.73	113.50

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	1020	DC	Sidechain
1	C	1021	DA	Sidechain
2	D	2002	DT	Sidechain
2	D	2005	DA	Sidechain
2	D	2007	DG	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	C	451	0	245	24	0
2	D	514	0	284	7	0
3	A	5537	0	5399	157	0
4	B	798	0	794	34	0
5	A	3	0	0	0	0
6	A	5	0	0	1	0
7	A	27	0	12	3	0
8	A	12	0	13	0	0
9	A	13	0	18	1	0
10	A	319	0	0	11	0
10	B	19	0	0	1	0
10	C	16	0	0	0	0
10	D	27	0	0	0	0
All	All	7741	0	6765	209	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 15.

All (209) 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)
3:A:392:ILE:HG12	10:A:4163:HOH:O	1.66	0.95
3:A:318:ARG:O	3:A:319:GLU:HB2	1.71	0.90
1:C:1020:DC:H5'	1:C:1020:DC:H6	1.39	0.85
3:A:273:MET:HE3	3:A:282:PRO:HB3	1.62	0.82
3:A:666:ALA:HB3	3:A:696:MET:HE2	1.61	0.81
4:B:41:ILE:O	4:B:45:ILE:HG12	1.82	0.79
3:A:599:ARG:HD3	3:A:620:LEU:HD11	1.62	0.79
3:A:534:ASP:CG	3:A:549:LYS:HG2	2.07	0.78
3:A:297:ILE:HD12	3:A:297:ILE:H	1.47	0.78
3:A:580:VAL:HG12	3:A:581:ALA:H	1.48	0.78
3:A:480:GLU:CD	7:A:1023:DCT:H2''	2.09	0.78
3:A:189:LYS:HG2	3:A:194:PRO:HD3	1.68	0.76
4:B:39:ALA:HB3	4:B:40:PRO:HD3	1.68	0.74
1:C:1006:DA:H2''	1:C:1007:DC:C5'	2.17	0.74
2:D:2004:DG:H5''	3:A:532:ALA:HA	1.71	0.72
3:A:194:PRO:HG2	3:A:195:GLU:OE2	1.89	0.71
3:A:645:ASP:HB3	3:A:665:ILE:HD13	1.74	0.70
4:B:100:LYS:HB3	4:B:100:LYS:NZ	2.07	0.69
1:C:1020:DC:H2''	1:C:1021:DA:H5'	1.74	0.69
1:C:1006:DA:H2''	1:C:1007:DC:H5'	1.75	0.69
3:A:273:MET:HG3	3:A:288:ARG:NH2	2.09	0.68
4:B:96:LYS:O	4:B:100:LYS:HG3	1.94	0.68
4:B:2:ASP:N	4:B:2:ASP:OD1	2.27	0.68
3:A:297:ILE:HD12	3:A:297:ILE:N	2.08	0.68
3:A:301:PRO:HA	3:A:306:GLN:HE21	1.58	0.68
3:A:302:LYS:H	3:A:306:GLN:NE2	1.93	0.66
3:A:492:ASP:O	3:A:495:GLU:HG3	1.96	0.65
3:A:580:VAL:HG12	3:A:581:ALA:N	2.12	0.65
3:A:545:LYS:O	3:A:549:LYS:HG3	1.97	0.64
1:C:1007:DC:H2''	1:C:1008:DG:C8	2.33	0.63
4:B:4:ILE:HG23	4:B:55:VAL:O	1.99	0.63
3:A:94:LEU:HB3	3:A:185:LEU:HD13	1.80	0.62
3:A:343:GLN:HG3	3:A:362:PRO:HG3	1.81	0.62
4:B:100:LYS:HB3	4:B:100:LYS:HZ2	1.63	0.62
4:B:13:ASP:HB3	10:B:4044:HOH:O	2.00	0.62
3:A:49:GLU:OE2	3:A:54:GLY:HA3	1.99	0.61
3:A:534:ASP:OD2	3:A:549:LYS:HG2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:317:THR:HG22	3:A:317:THR:O	2.00	0.61
3:A:632:MET:HA	3:A:635:GLU:HG2	1.81	0.61
1:C:1005:DA:H4'	3:A:294:VAL:HG21	1.83	0.61
3:A:537:ILE:O	3:A:541:VAL:HG22	2.01	0.60
3:A:570:GLN:HE22	3:A:606:PRO:HB3	1.65	0.60
3:A:142:ASP:HA	3:A:145:ARG:NH2	2.15	0.60
3:A:297:ILE:H	3:A:297:ILE:CD1	2.14	0.60
3:A:273:MET:HA	3:A:288:ARG:HE	1.66	0.60
3:A:195:GLU:H	3:A:195:GLU:CD	2.09	0.60
3:A:343:GLN:HG3	3:A:362:PRO:CG	2.30	0.60
3:A:628:LYS:HE3	3:A:679:TRP:CE2	2.37	0.59
3:A:553:LYS:O	3:A:557:GLU:HB2	2.02	0.59
3:A:155:VAL:O	3:A:156:ASP:C	2.45	0.59
3:A:339:ARG:HG2	3:A:339:ARG:HH11	1.68	0.59
1:C:1016:DG:H2''	1:C:1017:DT:H5'	1.85	0.59
2:D:2005:DA:H2'	2:D:2006:DT:H72	1.85	0.58
3:A:421:ASN:HB3	3:A:431:THR:OG1	2.02	0.58
3:A:276:HIS:CE1	3:A:278:ARG:HD2	2.39	0.58
4:B:82:LYS:O	4:B:83:ASN:HB2	2.04	0.58
2:D:2008:DG:H4'	3:A:432:HIS:O	2.04	0.58
3:A:394:LYS:HE3	10:A:4154:HOH:O	2.04	0.58
3:A:478:GLY:O	3:A:482:ARG:HG3	2.03	0.58
3:A:249:ARG:HG3	3:A:392:ILE:HD13	1.85	0.57
3:A:597:ASP:OD1	3:A:599:ARG:HD2	2.03	0.57
3:A:150:GLN:HB2	3:A:152:GLU:OE2	2.05	0.57
3:A:515:LEU:HD12	3:A:521:ALA:HA	1.87	0.57
3:A:195:GLU:CD	3:A:195:GLU:N	2.63	0.57
3:A:344:LYS:O	3:A:348:GLU:HG3	2.03	0.57
3:A:49:GLU:CD	3:A:52:ARG:HH12	2.13	0.56
3:A:91:THR:HB	3:A:181:LEU:HD13	1.88	0.55
3:A:326:TYR:HB3	4:B:92:GLY:HA2	1.88	0.55
3:A:85:ARG:HG3	3:A:222:TRP:CG	2.42	0.54
3:A:534:ASP:OD1	3:A:552:LYS:HE2	2.08	0.54
3:A:35:ARG:HB3	3:A:36:PRO:HD2	1.90	0.54
3:A:276:HIS:HD2	3:A:277:PRO:HD2	1.73	0.54
1:C:1021:DA:H5''	3:A:440:ILE:O	2.08	0.53
3:A:155:VAL:HG12	3:A:156:ASP:N	2.23	0.53
1:C:1006:DA:H1'	1:C:1007:DC:H5''	1.91	0.53
3:A:173:GLN:HG2	10:A:4430:HOH:O	2.08	0.53
3:A:370:GLU:OE2	3:A:387:LYS:HE3	2.07	0.53
3:A:237:LYS:HE3	3:A:237:LYS:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:404:LYS:HA	3:A:409:TYR:HE2	1.73	0.52
1:C:1013:DC:H3'	3:A:111:ARG:NH1	2.24	0.52
3:A:299:LYS:HD2	3:A:314:GLU:HB3	1.91	0.52
3:A:522:LYS:HE3	7:A:1023:DCT:O1A	2.09	0.52
4:B:95:SER:OG	4:B:98:GLN:HG3	2.09	0.52
3:A:296:GLY:O	3:A:316:ASP:HB3	2.10	0.51
3:A:480:GLU:OE2	7:A:1023:DCT:H2''	2.10	0.51
4:B:32:CYS:SG	4:B:34:PRO:HD2	2.50	0.51
3:A:456:GLY:HA2	3:A:471:GLN:OE1	2.11	0.51
3:A:273:MET:HG3	3:A:288:ARG:HH21	1.76	0.51
3:A:546:GLU:CD	3:A:546:GLU:H	2.18	0.51
3:A:301:PRO:HA	3:A:306:GLN:NE2	2.23	0.51
3:A:339:ARG:HG2	3:A:339:ARG:NH1	2.25	0.51
3:A:412:GLU:HB3	10:A:4102:HOH:O	2.10	0.51
3:A:440:ILE:O	3:A:452:ARG:NH1	2.44	0.51
3:A:158:MET:HA	3:A:161:TRP:CD2	2.45	0.50
3:A:440:ILE:HD12	3:A:700:TRP:CZ3	2.46	0.50
4:B:49:TYR:HB3	4:B:53:LEU:HD22	1.93	0.50
4:B:70:TYR:HB2	4:B:72:ILE:CD1	2.41	0.50
3:A:284:PRO:HA	3:A:288:ARG:HH22	1.77	0.50
1:C:1013:DC:C2'	3:A:111:ARG:HH12	2.25	0.50
3:A:182:LEU:O	3:A:186:LEU:HG	2.12	0.50
2:D:2005:DA:H2'	2:D:2006:DT:C7	2.41	0.50
3:A:35:ARG:HB3	3:A:36:PRO:CD	2.42	0.50
1:C:1009:DA:H2''	1:C:1010:DC:O5'	2.12	0.49
3:A:452:ARG:HG3	3:A:700:TRP:HB3	1.93	0.49
3:A:319:GLU:O	4:B:90:LYS:HA	2.13	0.49
3:A:152:GLU:CD	3:A:152:GLU:H	2.21	0.49
4:B:22:ALA:HB3	4:B:53:LEU:HB2	1.95	0.49
3:A:64:TYR:CZ	9:A:4009:PG4:H21	2.48	0.49
1:C:1006:DA:H2''	1:C:1007:DC:H5''	1.93	0.49
1:C:1021:DA:O4'	3:A:439:GLN:HA	2.13	0.49
3:A:112:SER:HA	3:A:115:LEU:HD12	1.94	0.48
1:C:1006:DA:C2'	1:C:1007:DC:H5''	2.42	0.48
3:A:336:PRO:HB2	3:A:389:TYR:CD1	2.49	0.48
3:A:342:ILE:O	3:A:346:LEU:HG	2.13	0.48
3:A:282:PRO:O	3:A:284:PRO:HD3	2.14	0.48
3:A:264:TRP:CZ3	3:A:345:LYS:HE2	2.49	0.48
3:A:58:PHE:HB2	10:A:4241:HOH:O	2.13	0.47
3:A:631:GLU:HG2	10:A:4297:HOH:O	2.12	0.47
3:A:35:ARG:HG3	10:A:4337:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1020:DC:H5'	1:C:1020:DC:C6	2.32	0.47
4:B:5:ILE:HG22	4:B:7:LEU:HD23	1.96	0.47
3:A:130:TYR:O	3:A:134:GLU:HG2	2.13	0.47
4:B:2:ASP:O	4:B:3:LYS:HD3	2.15	0.47
3:A:236:THR:O	3:A:240:GLU:HB2	2.15	0.46
3:A:420:VAL:O	3:A:422:PRO:HD3	2.14	0.46
4:B:19:ALA:CB	4:B:23:ILE:HD11	2.45	0.46
3:A:305:ALA:HA	3:A:310:ARG:NH1	2.31	0.46
3:A:515:LEU:HD13	3:A:520:ASN:OD1	2.15	0.46
3:A:580:VAL:CG1	3:A:581:ALA:H	2.17	0.46
3:A:249:ARG:HG3	3:A:392:ILE:CD1	2.46	0.46
3:A:506:HIS:CG	3:A:522:LYS:HG3	2.51	0.46
3:A:336:PRO:HB2	3:A:389:TYR:CE1	2.50	0.46
3:A:35:ARG:NE	3:A:165:GLU:OE1	2.47	0.46
3:A:525:ILE:HG23	3:A:526:TYR:N	2.31	0.46
3:A:343:GLN:HG3	3:A:362:PRO:HG2	1.98	0.45
1:C:1003:DA:H2''	1:C:1004:DA:OP2	2.16	0.45
3:A:79:ARG:HG2	3:A:80:GLU:N	2.31	0.45
4:B:77:THR:CG2	4:B:79:LEU:HD13	2.45	0.45
3:A:347:GLN:C	3:A:349:ALA:H	2.25	0.45
1:C:1013:DC:H2'	3:A:111:ARG:HH12	1.81	0.45
3:A:58:PHE:O	3:A:90:ASP:HA	2.17	0.45
3:A:229:ARG:NH1	6:A:999:SO4:O4	2.49	0.45
1:C:1021:DA:H2'	1:C:1022:2DT:H72	1.99	0.45
4:B:37:MET:O	4:B:40:PRO:HD2	2.16	0.45
3:A:17:LYS:HD2	3:A:76:GLN:NE2	2.32	0.45
3:A:237:LYS:O	3:A:241:GLU:HG3	2.17	0.45
3:A:35:ARG:O	3:A:36:PRO:C	2.61	0.44
3:A:111:ARG:HH21	3:A:114:LYS:HD2	1.82	0.44
3:A:379:LYS:HE3	10:A:4162:HOH:O	2.17	0.44
1:C:1016:DG:H5''	3:A:359:LYS:HD3	1.98	0.44
4:B:12:PHE:CE2	4:B:66:THR:HG21	2.53	0.44
4:B:67:ALA:HB3	4:B:68:PRO:HD3	1.98	0.44
1:C:1016:DG:H5''	3:A:359:LYS:CD	2.48	0.44
3:A:493:ASN:ND2	10:A:4271:HOH:O	2.40	0.44
4:B:13:ASP:C	4:B:13:ASP:OD2	2.60	0.44
4:B:27:PHE:CE1	4:B:79:LEU:HD22	2.53	0.44
4:B:19:ALA:O	4:B:20:ASP:HB2	2.18	0.44
3:A:96:ARG:NH1	10:A:4105:HOH:O	2.48	0.44
1:C:1020:DC:H6	1:C:1020:DC:C5'	2.22	0.44
2:D:2025:DC:H2''	2:D:2026:DG:C8	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:267:PRO:HG3	4:B:31:TRP:CZ3	2.53	0.43
1:C:1020:DC:H2''	1:C:1021:DA:C5'	2.47	0.43
3:A:146:MET:O	3:A:150:GLN:HG2	2.19	0.43
3:A:13:GLU:H	3:A:13:GLU:CD	2.26	0.43
3:A:273:MET:HA	3:A:288:ARG:NE	2.33	0.43
3:A:233:PRO:HB2	3:A:456:GLY:O	2.18	0.43
3:A:443:VAL:HG23	3:A:444:ARG:HG3	2.01	0.43
3:A:484:LEU:O	3:A:488:MET:HG2	2.18	0.43
3:A:548:GLY:O	3:A:552:LYS:HG3	2.19	0.43
3:A:597:ASP:OD1	3:A:599:ARG:CD	2.67	0.43
2:D:2018:DT:H2''	2:D:2019:DC:O5'	2.18	0.42
3:A:574:VAL:O	3:A:589:LYS:HE3	2.19	0.42
1:C:1019:DC:C2	1:C:1020:DC:C5	3.07	0.42
3:A:310:ARG:O	3:A:310:ARG:HG2	2.18	0.42
3:A:106:ASP:HB2	3:A:110:LEU:HD22	2.01	0.42
3:A:267:PRO:HB3	3:A:287:PRO:HB3	2.02	0.42
3:A:633:LEU:HD21	3:A:669:VAL:HA	2.01	0.42
3:A:474:ILE:HD12	3:A:474:ILE:N	2.34	0.42
2:D:2019:DC:H2''	2:D:2020:DG:C8	2.55	0.42
3:A:58:PHE:CB	3:A:61:GLY:HA3	2.50	0.42
3:A:308:GLU:OE2	3:A:310:ARG:NH2	2.52	0.42
3:A:545:LYS:HG3	3:A:546:GLU:OE2	2.20	0.42
3:A:24:TYR:OH	3:A:29:ALA:HA	2.20	0.42
3:A:309:GLY:HA2	10:A:4281:HOH:O	2.19	0.42
4:B:25:VAL:HA	4:B:56:ALA:O	2.20	0.42
4:B:77:THR:HG22	4:B:79:LEU:HD13	2.00	0.42
3:A:253:LEU:CD1	3:A:392:ILE:HD12	2.50	0.41
3:A:135:MET:CE	3:A:136:LYS:HG3	2.50	0.41
3:A:279:THR:OG1	3:A:280:GLY:N	2.53	0.41
4:B:83:ASN:C	4:B:85:GLU:H	2.28	0.41
3:A:569:ILE:HG21	3:A:609:ALA:HB1	2.02	0.41
3:A:570:GLN:NE2	3:A:606:PRO:HB3	2.34	0.41
3:A:145:ARG:HB2	3:A:145:ARG:NH1	2.36	0.41
3:A:145:ARG:HB2	3:A:145:ARG:HH11	1.86	0.41
3:A:432:HIS:ND1	3:A:432:HIS:N	2.68	0.41
4:B:99:LEU:O	4:B:102:PHE:HB3	2.21	0.41
3:A:111:ARG:NH2	3:A:114:LYS:HD2	2.36	0.41
3:A:235:ASP:HB2	3:A:459:HIS:CE1	2.56	0.41
3:A:639:LYS:HG3	3:A:644:GLY:HA3	2.03	0.41
3:A:276:HIS:HA	3:A:277:PRO:HD3	1.94	0.41
3:A:446:PRO:O	3:A:447:TYR:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:477:SER:O	3:A:478:GLY:C	2.65	0.40
3:A:211:LEU:O	3:A:212:GLU:C	2.65	0.40
3:A:250:ARG:HG3	3:A:389:TYR:CE2	2.56	0.40
3:A:343:GLN:O	3:A:347:GLN:HG3	2.21	0.40
4:B:22:ALA:HB3	4:B:53:LEU:CB	2.51	0.40
3:A:59:HIS:O	3:A:60:ASN:HB3	2.21	0.40
4:B:19:ALA:O	4:B:20:ASP:O	2.40	0.40
4:B:49:TYR:HB3	4:B:53:LEU:CD2	2.51	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	
3	A	696/698 (100%)	651 (94%)	36 (5%)	9 (1%)	9	12
4	B	104/108 (96%)	94 (90%)	8 (8%)	2 (2%)	6	7
All	All	800/806 (99%)	745 (93%)	44 (6%)	11 (1%)	9	11

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	156	ASP
3	A	319	GLU
3	A	318	ARG
3	A	580	VAL
4	B	83	ASN
4	B	20	ASP
3	A	278	ARG
3	A	279	THR
3	A	280	GLY
3	A	653	HIS

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Mol	Chain	Res	Type
3	A	579	TRP

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	
3	A	571/579 (99%)	551 (96%)	20 (4%)	32	48
4	B	83/87 (95%)	75 (90%)	8 (10%)	8	9
All	All	654/666 (98%)	626 (96%)	28 (4%)	26	39

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	78	ASN
3	A	110	LEU
3	A	159	GLU
3	A	176	VAL
3	A	181	LEU
3	A	189	LYS
3	A	237	LYS
3	A	297	ILE
3	A	314	GLU
3	A	332	VAL
3	A	343	GLN
3	A	410	VAL
3	A	426	VAL
3	A	499	GLU
3	A	536	LYS
3	A	551	LEU
3	A	557	GLU
3	A	599	ARG
3	A	624	LEU
3	A	686	PHE
4	B	2	ASP
4	B	14	THR

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Mol	Chain	Res	Type
4	B	53	LEU
4	B	54	THR
4	B	79	LEU
4	B	80	LEU
4	B	89	THR
4	B	99	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	227	GLN
3	A	276	HIS
3	A	303	ASN
3	A	306	GLN
3	A	343	GLN
3	A	347	GLN
3	A	450	GLN
3	A	539	GLN
3	A	570	GLN
4	B	62	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	2DT	C	1022	1,2	17,20,21	0.47	0	22,28,31	0.71	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
1	2DT	C	1022	1,2	-	0/7/18/19	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	1022	2DT	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 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
7	DCT	A	1023	5	27,28,28	0.69	0	37,43,43	1.05	3 (8%)
9	PG4	A	4009	-	12,12,12	1.03	1 (8%)	11,11,11	2.00	4 (36%)
6	SO4	A	999	-	4,4,4	0.93	0	6,6,6	0.20	0
8	MES	A	4008	-	12,12,12	0.91	0	15,16,16	0.74	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
9	PG4	A	4009	-	-	0/10/10/10	-
7	DCT	A	1023	5	-	2/22/31/31	0/2/2/2
8	MES	A	4008	-	-	1/6/14/14	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	4009	PG4	C4-C3	2.12	1.59	1.49

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	4009	PG4	C3-O2-C2	-4.07	95.44	113.26
7	A	1023	DCT	O3G-PG-O3B	-3.10	94.25	104.64
9	A	4009	PG4	C5-O3-C4	-2.91	100.55	113.26
9	A	4009	PG4	O2-C2-C1	-2.68	98.31	110.11
9	A	4009	PG4	O4-C6-C5	-2.39	99.47	110.35
7	A	1023	DCT	O2G-PG-O1G	2.29	119.77	110.83
7	A	1023	DCT	C2'-C1'-N1	2.07	116.32	112.40

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	4008	MES	C8-C7-N4-C5
7	A	1023	DCT	PA-O3A-PB-O3B
7	A	1023	DCT	PA-O3A-PB-O1B

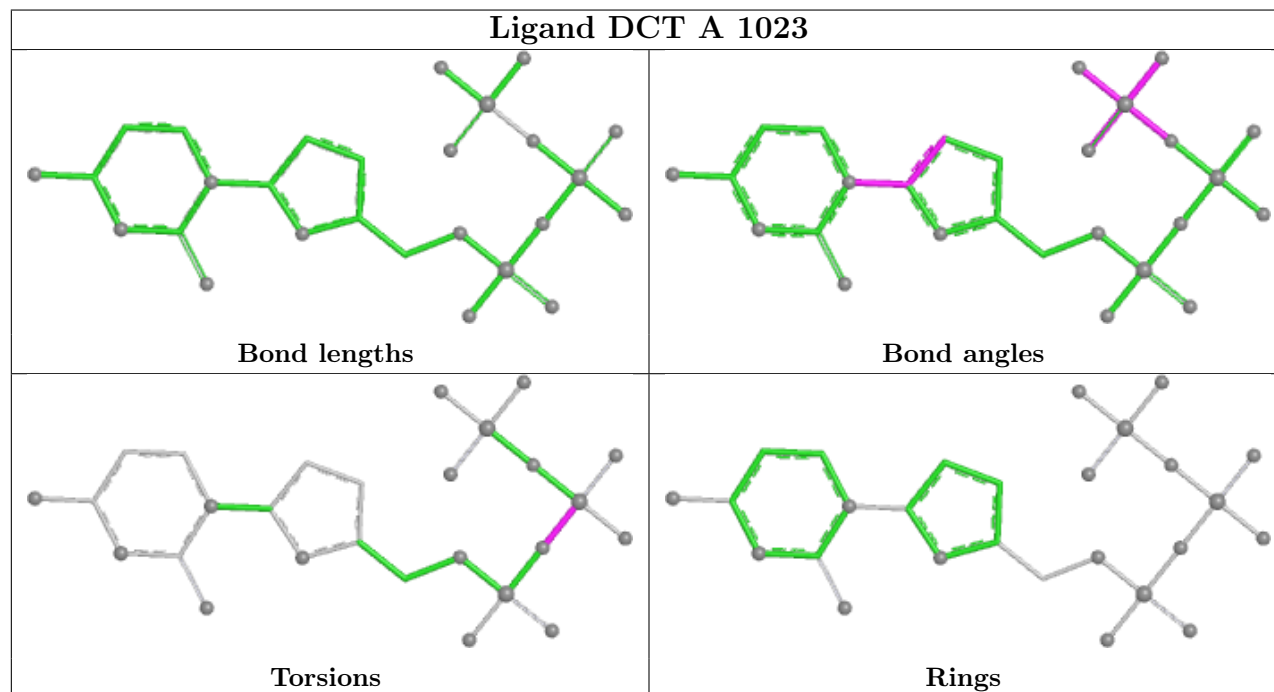
There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1023	DCT	3	0
9	A	4009	PG4	1	0
6	A	999	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 ⓘ

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	C	21/22 (95%)	0.98	5 (23%) 2 1	24, 83, 98, 100	0
2	D	25/26 (96%)	0.95	9 (36%) 1 0	19, 56, 100, 100	0
3	A	698/698 (100%)	-0.12	21 (3%) 52 54	13, 28, 68, 86	0
4	B	106/108 (98%)	0.54	6 (5%) 29 29	24, 47, 69, 73	0
All	All	850/854 (99%)	0.03	41 (4%) 35 35	13, 31, 74, 100	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	2024	DT	3.7
3	A	582	GLY	3.7
3	A	303	ASN	3.7
3	A	583	GLU	3.4
3	A	294	VAL	3.3
3	A	580	VAL	3.2
4	B	84	GLY	3.1
2	D	2025	DC	3.1
3	A	278	ARG	3.0
4	B	107	LEU	2.9
3	A	147	LEU	2.9
3	A	317	THR	2.8
1	C	1006	DA	2.8
3	A	13	GLU	2.7
3	A	581	ALA	2.6
1	C	1007	DC	2.5
4	B	48	GLU	2.5
1	C	1003	DA	2.5
4	B	99	LEU	2.5
2	D	2015	DC	2.4
2	D	2026	DG	2.4

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Mol	Chain	Res	Type	RSRZ
3	A	366	ASP	2.3
3	A	156	ASP	2.3
3	A	150	GLN	2.3
2	D	2019	DC	2.3
3	A	298	PHE	2.3
2	D	2018	DT	2.3
4	B	22	ALA	2.2
3	A	312	PRO	2.2
4	B	2	ASP	2.2
2	D	2023	DT	2.2
1	C	1002	DG	2.2
1	C	1008	DG	2.2
2	D	2020	DG	2.2
3	A	284	PRO	2.2
3	A	297	ILE	2.1
3	A	318	ARG	2.1
3	A	293	LYS	2.1
3	A	279	THR	2.0
3	A	292	PRO	2.0
2	D	2002	DT	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	2DT	C	1022	19/20	0.98	0.07	19,22,26,27	0

6.3 Carbohydrates [i](#)

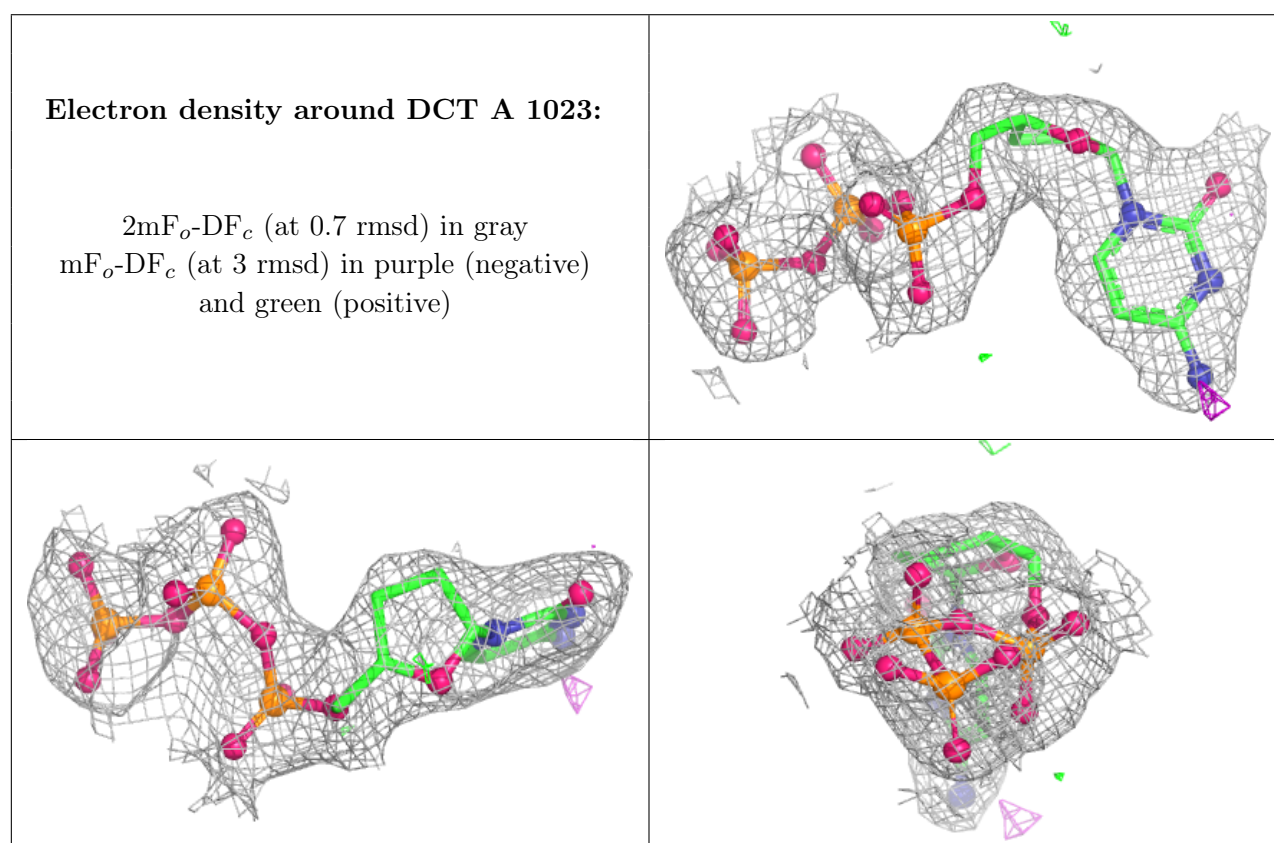
There are no oligosaccharides in this entry.

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
9	PG4	A	4009	13/13	0.87	0.17	56,59,63,64	0
8	MES	A	4008	12/12	0.92	0.12	40,42,46,46	0
5	MG	A	4003	1/1	0.95	0.07	22,22,22,22	0
5	MG	A	4002	1/1	0.96	0.06	31,31,31,31	0
6	SO4	A	999	5/5	0.96	0.10	43,44,45,46	0
7	DCT	A	1023	27/27	0.97	0.07	17,22,26,26	0
5	MG	A	4001	1/1	0.99	0.05	16,16,16,16	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.