



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:37 AM UTC

PDB ID : 6DX0 / pdb_00006dx0
Title : Hermes transposase deletion dimer complex with (A/T) DNA
Authors : Dyda, F.; Voth, A.R.; Hickman, A.B.
Deposited on : 2018-06-28
Resolution : 2.90 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

<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
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

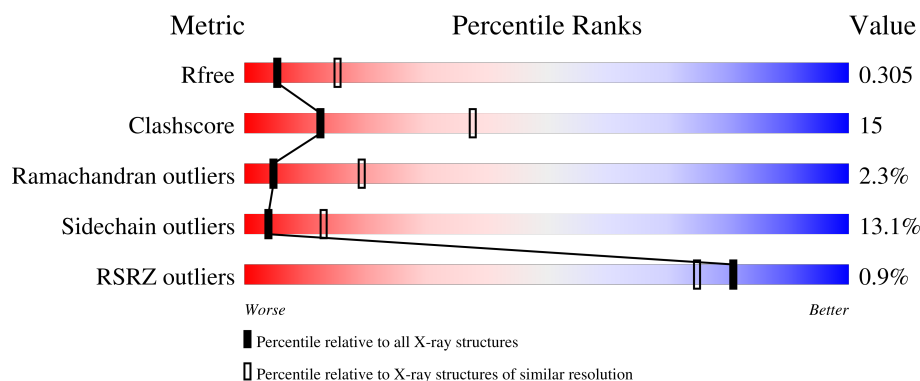
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 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	517	14% 56% 32% 5% 7%
2	B	16	19% 81%
3	C	25	48% 48% .
4	D	7	100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4867 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 Hermes transposase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	483	Total	C	N	O	S	0	0	0
			3888	2489	658	722	19			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	GLY	-	expression tag	UNP Q25438
A	77	SER	-	expression tag	UNP Q25438
A	78	HIS	-	expression tag	UNP Q25438
A	79	MET	-	expression tag	UNP Q25438
A	128	GLY	LYS	conflict	UNP Q25438
A	?	-	ASP	deletion	UNP Q25438
A	?	-	ILE	deletion	UNP Q25438
A	?	-	SER	deletion	UNP Q25438
A	?	-	THR	deletion	UNP Q25438
A	?	-	THR	deletion	UNP Q25438
A	?	-	SER	deletion	UNP Q25438
A	?	-	PHE	deletion	UNP Q25438
A	?	-	PHE	deletion	UNP Q25438
A	?	-	PHE	deletion	UNP Q25438
A	?	-	PRO	deletion	UNP Q25438
A	?	-	GLN	deletion	UNP Q25438
A	?	-	LEU	deletion	UNP Q25438
A	?	-	THR	deletion	UNP Q25438
A	?	-	GLN	deletion	UNP Q25438
A	?	-	ASN	deletion	UNP Q25438
A	?	-	ASN	deletion	UNP Q25438
A	?	-	SER	deletion	UNP Q25438
A	?	-	ARG	deletion	UNP Q25438
A	?	-	GLU	deletion	UNP Q25438
A	?	-	PRO	deletion	UNP Q25438
A	519	SER	CYS	engineered mutation	UNP Q25438

- Molecule 2 is a DNA chain called DNA (5'-D(*AP*GP*AP*GP*AP*AP*CP*AP*AP*CP*AP*AP*CP*AP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	16	Total	C	N	O	P	0	0	0
			330	157	74	84	15			

- Molecule 3 is a DNA chain called DNA (25-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	25	Total	C	N	O	P	0	0	0
			504	244	77	159	24			

- Molecule 4 is a DNA chain called DNA (5'-D(*GP*CP*GP*TP*GP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	7	Total	C	N	O	P	0	0	0
			144	69	30	39	6			

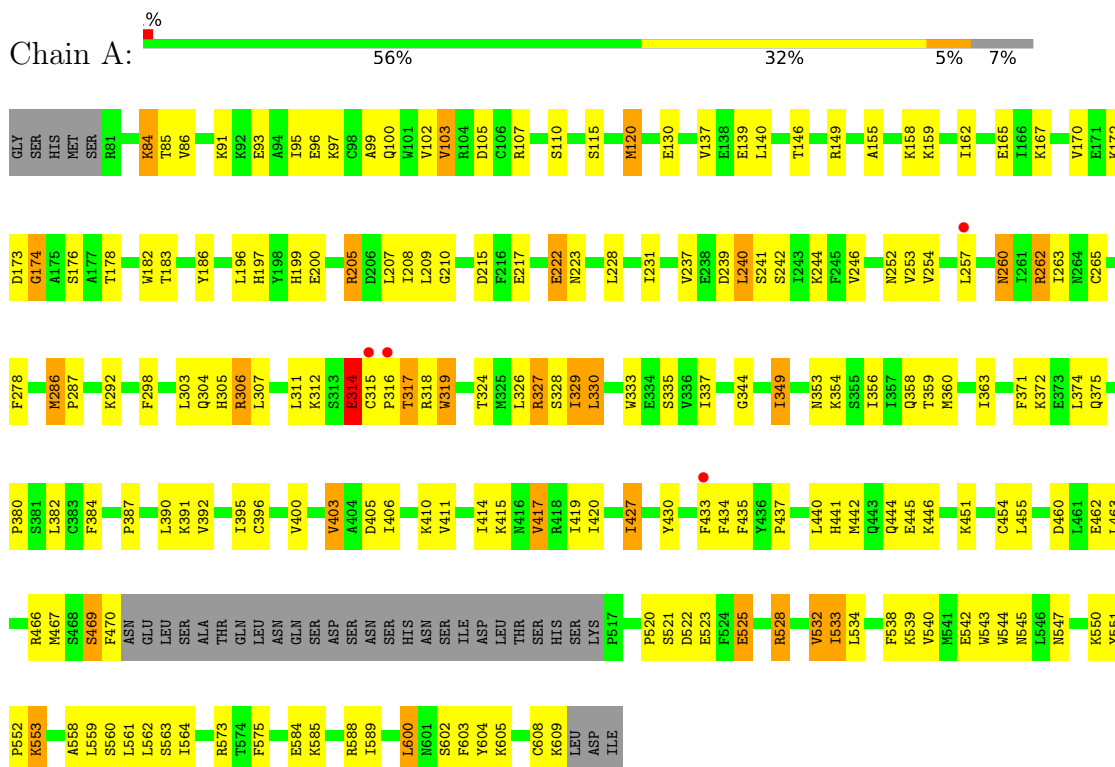
- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Na	0	0
			1	1		

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.

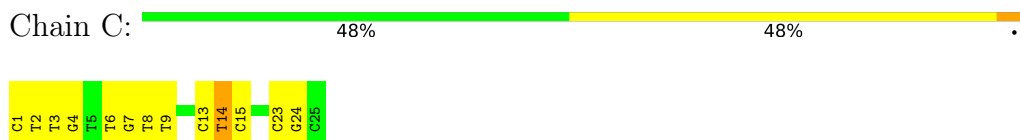
• Molecule 1: Hermes transposase



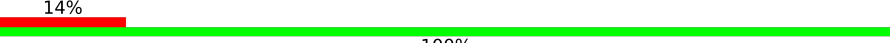
• Molecule 2: DNA (5'-D(*AP*GP*AP*GP*AP*AP*CP*AP*AP*CP*AP*AP*CP*AP*AP*G)-3')



• Molecule 3: DNA (25-MER)



- Molecule 4: DNA (5'-D(*GP*CP*GP*TP*GP*AP*A)-3')

Chain D:  14% 100%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	121.60Å 135.50Å 102.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.44 – 2.90 28.44 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.6 (28.44-2.90) 90.5 (28.44-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.225 , 0.305 0.228 , 0.305	Depositor DCC
R_{free} test set	869 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	79.8	Xtriage
Anisotropy	0.909	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4867	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.28	0/3962	0.54	1/5343 (0.0%)
2	B	0.42	0/374	0.64	0/575
3	C	0.46	0/560	0.75	1/863 (0.1%)
4	D	0.40	0/162	0.60	0/249
All	All	0.32	0/5058	0.58	2/7030 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	14	DT	O5'-C5'-C4'	5.11	118.46	110.80
1	A	344	GLY	N-CA-C	-5.05	108.64	115.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3888	0	3957	102	0
2	B	330	0	178	21	0
3	C	504	0	289	14	0
4	D	144	0	79	0	0
5	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4867	0	4503	132	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 (132) 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)
2:B:7:DC:H2''	2:B:8:DA:C8	2.12	0.84
1:A:372:LYS:NZ	2:B:1:DA:N7	2.29	0.79
1:A:444:GLN:O	1:A:446:LYS:N	2.17	0.78
1:A:573:ARG:HD3	2:B:2:DG:H1'	1.65	0.78
2:B:4:DG:H2''	2:B:5:DA:H5''	1.68	0.76
3:C:6:DT:H2''	3:C:7:DG:H5'	1.68	0.75
2:B:10:DC:H2'	2:B:11:DA:C8	2.22	0.74
1:A:330:LEU:HD11	1:A:358:GLN:HG2	1.69	0.74
3:C:3:DT:H2'	3:C:4:DG:C8	2.23	0.73
3:C:6:DT:H2'	3:C:7:DG:C8	2.24	0.73
1:A:304:GLN:OE1	1:A:311:LEU:N	2.25	0.69
1:A:196:LEU:HD22	1:A:209:LEU:HD11	1.75	0.68
1:A:430:TYR:HB3	1:A:442:MET:HE1	1.73	0.68
1:A:210:GLY:HA3	1:A:231:ILE:HD12	1.76	0.68
1:A:437:PRO:HB2	1:A:532:VAL:HG11	1.75	0.68
1:A:200:GLU:OE1	1:A:205:ARG:NH2	2.27	0.68
2:B:2:DG:H2'	2:B:3:DA:H8	1.59	0.68
1:A:318:ARG:NH2	3:C:15:DC:OP2	2.24	0.68
1:A:375:GLN:O	1:A:573:ARG:NH2	2.27	0.67
3:C:3:DT:H2''	3:C:4:DG:H5'	1.75	0.67
1:A:170:VAL:HG21	1:A:237:VAL:HG13	1.77	0.66
3:C:14:DT:H6	3:C:14:DT:H5''	1.60	0.66
3:C:1:DC:H2'	3:C:2:DT:C6	2.34	0.63
1:A:419:ILE:HG13	1:A:420:ILE:HG12	1.81	0.62
1:A:396:CYS:HA	1:A:410:LYS:HD3	1.81	0.61
2:B:4:DG:C2'	2:B:5:DA:H5''	2.30	0.61
2:B:2:DG:H2'	2:B:3:DA:C8	2.36	0.60
1:A:356:ILE:HG23	1:A:406:ILE:HD11	1.83	0.60
1:A:208:ILE:HD12	1:A:602:SER:HB2	1.83	0.60
1:A:437:PRO:HG2	1:A:538:PHE:HZ	1.66	0.59
2:B:9:DA:N6	3:C:8:DT:O4	2.17	0.59
1:A:292:LYS:HE3	1:A:314:GLU:HB2	1.86	0.58
1:A:403:VAL:HG23	1:A:405:ASP:H	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:6:DA:C2'	2:B:7:DC:H5''	2.35	0.56
1:A:246:VAL:HG22	1:A:263:ILE:HB	1.86	0.56
1:A:304:GLN:HA	1:A:307:LEU:HD23	1.88	0.56
1:A:437:PRO:HG2	1:A:538:PHE:CZ	2.40	0.56
1:A:239:ASP:O	1:A:241:SER:N	2.39	0.56
1:A:380:PRO:HB3	1:A:562:LEU:HB3	1.88	0.56
1:A:172:LYS:O	1:A:174:GLY:N	2.38	0.54
1:A:312:LYS:HG2	1:A:328:SER:CB	2.37	0.54
1:A:312:LYS:HG2	1:A:328:SER:HB3	1.89	0.54
1:A:523:GLU:HG3	1:A:551:TYR:HD2	1.71	0.54
1:A:178:THR:HG22	1:A:246:VAL:HB	1.90	0.53
1:A:542:GLU:HA	1:A:545:ASN:HD22	1.74	0.53
1:A:182:TRP:CD1	1:A:575:PHE:HB3	2.44	0.53
1:A:411:VAL:O	1:A:414:ILE:HG13	2.09	0.53
2:B:4:DG:C3'	2:B:5:DA:H5''	2.40	0.52
1:A:252:ASN:OD1	1:A:253:VAL:N	2.43	0.52
1:A:105:ASP:O	1:A:107:ARG:HG2	2.09	0.52
2:B:7:DC:H2''	2:B:8:DA:H8	1.70	0.52
1:A:158:LYS:HD2	1:A:603:PHE:CE2	2.45	0.52
1:A:305:HIS:CD2	1:A:306:ARG:HG2	2.45	0.52
1:A:333:TRP:O	1:A:337:ILE:HG23	2.10	0.52
1:A:96:GLU:O	1:A:100:GLN:HG3	2.10	0.52
2:B:7:DC:H5''	2:B:7:DC:H6	1.74	0.52
1:A:435:PHE:CD1	1:A:558:ALA:HA	2.44	0.51
1:A:260:ASN:N	1:A:260:ASN:OD1	2.42	0.51
1:A:315:CYS:C	1:A:317:THR:H	2.18	0.51
1:A:552:PRO:HD2	1:A:553:LYS:HE3	1.93	0.51
1:A:538:PHE:HE1	1:A:543:TRP:HB2	1.76	0.51
3:C:13:DC:H2'	3:C:14:DT:H5''	1.93	0.50
1:A:540:VAL:O	1:A:544:TRP:HD1	1.93	0.50
1:A:463:LEU:O	1:A:467:MET:HG2	2.12	0.49
2:B:8:DA:H8	2:B:8:DA:OP2	1.96	0.49
1:A:349:ILE:HG12	1:A:349:ILE:O	2.13	0.49
3:C:23:DC:H2''	3:C:24:DG:C8	2.48	0.49
1:A:298:PHE:CE1	1:A:307:LEU:HD21	2.47	0.49
1:A:382:LEU:HD22	1:A:564:ILE:HB	1.95	0.49
1:A:327:ARG:HB3	1:A:327:ARG:HH11	1.78	0.49
1:A:298:PHE:HE1	1:A:307:LEU:HD21	1.77	0.49
1:A:110:SER:OG	2:B:5:DA:H5'	2.13	0.48
1:A:520:PRO:HA	1:A:523:GLU:HB3	1.94	0.48
1:A:155:ALA:O	1:A:159:LYS:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:TRP:CH2	1:A:354:LYS:HB2	2.48	0.48
1:A:359:THR:O	1:A:363:ILE:HG12	2.14	0.48
1:A:600:LEU:O	1:A:603:PHE:N	2.46	0.48
1:A:199:HIS:CE1	1:A:559:LEU:HD13	2.48	0.48
2:B:7:DC:H4'	2:B:7:DC:OP1	2.14	0.47
1:A:525:GLU:OE2	1:A:528:ARG:NH2	2.45	0.47
1:A:207:LEU:HD12	1:A:602:SER:OG	2.15	0.47
1:A:286:MET:HB3	1:A:287:PRO:HD3	1.96	0.47
1:A:560:SER:O	1:A:563:SER:OG	2.29	0.47
2:B:15:DA:H2''	2:B:16:DG:C8	2.50	0.47
1:A:315:CYS:N	1:A:316:PRO:HD2	2.29	0.47
1:A:146:THR:O	1:A:149:ARG:N	2.48	0.47
1:A:99:ALA:O	1:A:103:VAL:HG13	2.15	0.47
1:A:550:LYS:O	1:A:552:PRO:HD3	2.14	0.46
2:B:6:DA:H2''	2:B:7:DC:O4'	2.16	0.46
1:A:217:GLU:HB3	1:A:223:ASN:ND2	2.31	0.46
1:A:84:LYS:H	1:A:84:LYS:HG3	1.46	0.46
1:A:427:ILE:HA	1:A:430:TYR:CD1	2.51	0.45
1:A:120:MET:HE2	1:A:120:MET:HB3	1.72	0.45
3:C:2:DT:H2'	3:C:3:DT:H71	1.98	0.45
3:C:3:DT:H2'	3:C:4:DG:H8	1.76	0.45
1:A:523:GLU:HG3	1:A:551:TYR:CD2	2.51	0.45
1:A:539:LYS:O	1:A:542:GLU:HG2	2.17	0.44
1:A:588:ARG:HA	1:A:588:ARG:HD2	1.75	0.44
1:A:167:LYS:HD2	1:A:167:LYS:HA	1.67	0.44
1:A:337:ILE:HG22	1:A:349:ILE:CD1	2.46	0.44
1:A:93:GLU:HA	1:A:96:GLU:HB3	2.00	0.44
1:A:315:CYS:N	1:A:316:PRO:CD	2.81	0.44
1:A:371:PHE:O	1:A:375:GLN:HG3	2.18	0.44
1:A:391:LYS:O	1:A:395:ILE:HG13	2.17	0.43
1:A:451:LYS:O	1:A:455:LEU:HG	2.18	0.43
3:C:9:DT:C6	3:C:9:DT:H5'	2.54	0.43
1:A:382:LEU:HB2	1:A:564:ILE:O	2.19	0.43
1:A:538:PHE:CE1	1:A:543:TRP:HB2	2.53	0.43
3:C:1:DC:H2''	3:C:2:DT:H5'	2.01	0.43
1:A:311:LEU:HD22	1:A:329:ILE:HA	2.00	0.43
2:B:2:DG:H2''	2:B:3:DA:O4'	2.19	0.42
1:A:278:PHE:HE1	1:A:360:MET:HE3	1.84	0.42
1:A:257:LEU:HB2	1:A:262:ARG:HG3	2.01	0.42
1:A:186:TYR:OH	1:A:584:GLU:HG3	2.20	0.42
1:A:603:PHE:C	1:A:605:LYS:H	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:DA:H2''	2:B:2:DG:O4'	2.20	0.42
1:A:318:ARG:O	1:A:319:TRP:C	2.63	0.41
1:A:600:LEU:HD12	1:A:600:LEU:HA	1.84	0.41
1:A:85:THR:HG22	1:A:86:VAL:H	1.84	0.41
1:A:415:LYS:HE3	1:A:415:LYS:HB2	1.78	0.41
1:A:469:SER:HG	1:A:470:PHE:HD2	1.68	0.41
1:A:176:SER:OG	1:A:197:HIS:HB2	2.20	0.41
1:A:392:VAL:HG11	1:A:417:VAL:HG21	2.03	0.41
1:A:433:PHE:HD2	1:A:442:MET:SD	2.43	0.41
2:B:1:DA:C6	2:B:2:DG:C6	3.09	0.41
1:A:533:ILE:HD12	1:A:533:ILE:HA	1.79	0.41
1:A:91:LYS:HB3	1:A:91:LYS:HE2	1.63	0.41
1:A:434:PHE:HD2	1:A:435:PHE:CD2	2.38	0.40
1:A:222:GLU:H	1:A:222:GLU:HG3	1.66	0.40
1:A:384:PHE:C	1:A:387:PRO:HD2	2.47	0.40
1:A:462:GLU:O	1:A:466:ARG:HG2	2.22	0.40
1:A:162:ILE:HA	1:A:165:GLU:OE1	2.21	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
1	A	479/517 (93%)	421 (88%)	47 (10%)	11 (2%)	5 19

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	LEU
1	A	314	GLU
1	A	445	GLU
1	A	319	TRP

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Mol	Chain	Res	Type
1	A	173	ASP
1	A	215	ASP
1	A	604	TYR
1	A	353	ASN
1	A	469	SER
1	A	441	HIS
1	A	174	GLY

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	442/475 (93%)	384 (87%)	58 (13%)	4 13

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	LYS
1	A	95	ILE
1	A	97	LYS
1	A	102	VAL
1	A	103	VAL
1	A	115	SER
1	A	120	MET
1	A	130	GLU
1	A	137	VAL
1	A	139	GLU
1	A	140	LEU
1	A	183	THR
1	A	205	ARG
1	A	222	GLU
1	A	228	LEU
1	A	240	LEU
1	A	242	SER
1	A	244	LYS
1	A	254	VAL

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Mol	Chain	Res	Type
1	A	260	ASN
1	A	262	ARG
1	A	265	CYS
1	A	286	MET
1	A	303	LEU
1	A	306	ARG
1	A	314	GLU
1	A	317	THR
1	A	324	THR
1	A	326	LEU
1	A	327	ARG
1	A	329	ILE
1	A	330	LEU
1	A	335	SER
1	A	349	ILE
1	A	374	LEU
1	A	390	LEU
1	A	400	VAL
1	A	403	VAL
1	A	417	VAL
1	A	427	ILE
1	A	440	LEU
1	A	454	CYS
1	A	460	ASP
1	A	521	SER
1	A	522	ASP
1	A	525	GLU
1	A	528	ARG
1	A	532	VAL
1	A	533	ILE
1	A	534	LEU
1	A	547	ASN
1	A	553	LYS
1	A	561	LEU
1	A	585	LYS
1	A	589	ILE
1	A	600	LEU
1	A	608	CYS
1	A	609	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	134	HIS
1	A	234	GLN
1	A	362	ASN
1	A	416	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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	483/517 (93%)	-0.00	4 (0%) 82 77	63, 101, 152, 183	0
2	B	16/16 (100%)	-0.12	0 100 100	66, 93, 104, 108	0
3	C	25/25 (100%)	-0.02	0 100 100	69, 86, 107, 126	0
4	D	7/7 (100%)	0.06	1 (14%) 6 5	81, 87, 97, 102	0
All	All	531/565 (93%)	-0.01	5 (0%) 81 75	63, 99, 150, 183	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	1	DG	2.9
1	A	315	CYS	2.9
1	A	257	LEU	2.9
1	A	316	PRO	2.8
1	A	433	PHE	2.3

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
5	NA	A	701	1/1	0.94	0.20	71,71,71,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.