



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 15, 2018 – 12:00 PM EST

PDB ID : 3LDS
Title : Crystal structure of RB69 gp43 with DNA and dATP opposite 8-oxoG
Authors : Hogg, M.; Midkiff, J.; Rudnicki, J.; Reha-Krantz, L.; Doublié, S.; Wallace, S.S.
Deposited on : 2010-01-13
Resolution : 3.00 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

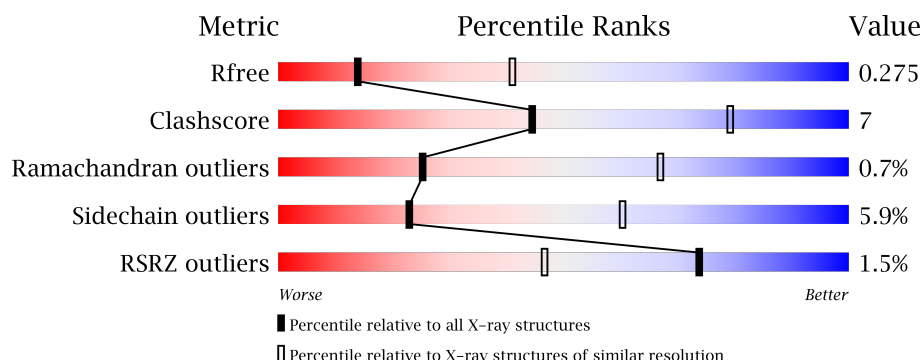
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (3.00-3.00)

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.

Mol	Chain	Length	Quality of chain
1	A	903	<div> <div>7%</div> <div>79%</div> <div>20%</div> <div>.</div> </div>
2	P	14	<div> <div>7%</div> <div>71%</div> <div>29%</div> </div>
3	T	18	<div> <div>61%</div> <div>39%</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8077 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 DNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	903	Total	C	N	O	S	0	0	0
			7371	4734	1226	1378	33			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ALA	ASP	engineered	UNP Q38087
A	327	ALA	ASP	engineered	UNP Q38087
A	561	ALA	LEU	engineered	UNP Q38087

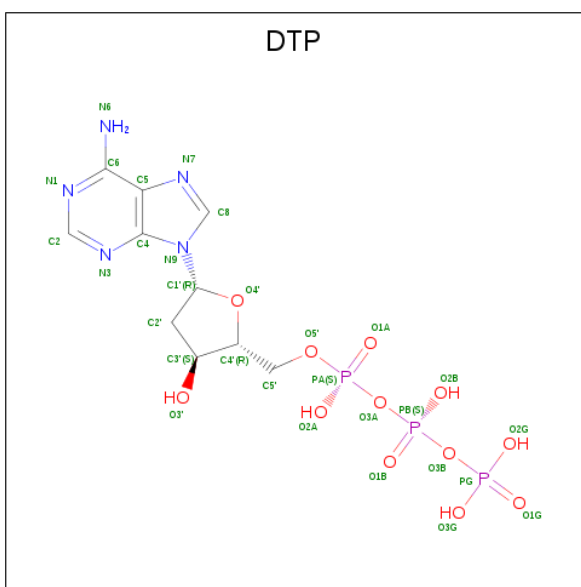
- Molecule 2 is a DNA chain called DNA (5'-D(*GP*CP*GP*GP*CP*TP*GP*TP*CP*AP*TP*AP*AP*(DDG))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	14	Total	C	N	O	P	0	0	0
			286	137	55	81	13			

- Molecule 3 is a DNA chain called DNA (5'-D(*CP*AP*(8OG)P*CP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*CP*GP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	18	Total	C	N	O	P	16	0	0
			366	174	69	106	17			

- Molecule 4 is 2'-DEOXYADENOSINE 5'-TRIPHOSPHATE (three-letter code: DTP) (formula: C₁₀H₁₆N₅O₁₂P₃).

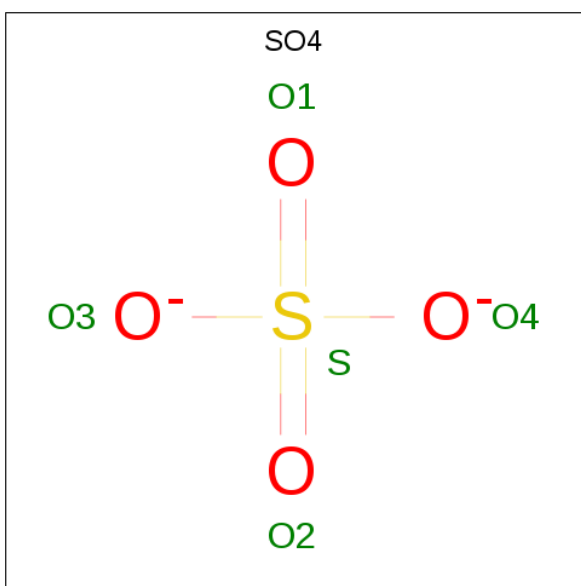


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			30	10	5	12	3		

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

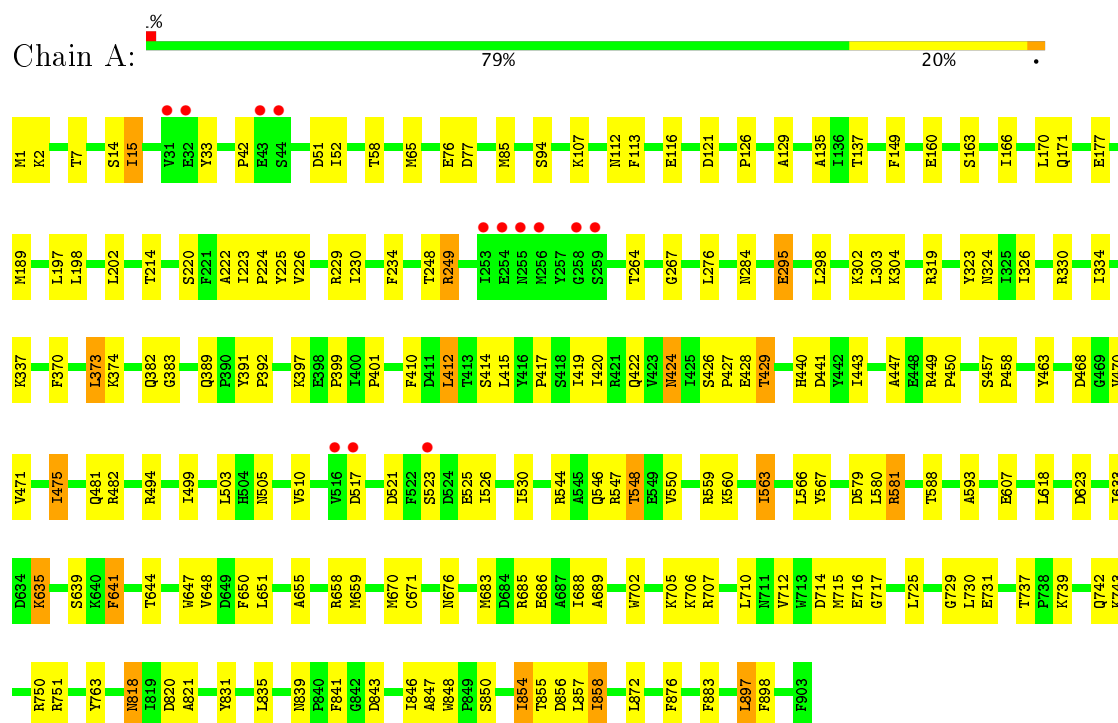
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	10	Total	O	0	0
			10	10		
7	P	1	Total	O	0	0
			1	1		

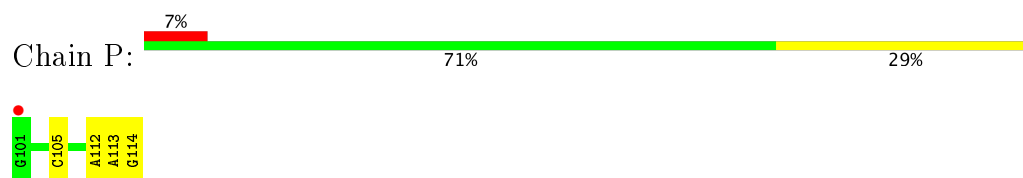
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: DNA polymerase



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



• Molecule 3: DNA (5'-D(*CP*AP*(8OG)P*CP*TP*TP*AP*TP*GP*AP*CP*AP*GP*CP*CP*GP*CP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.01Å 121.00Å 128.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.82 – 3.00 29.45 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.8 (29.82-3.00) 91.8 (29.45-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.6.0041	Depositor
R, R_{free}	0.242 , 0.295 0.208 , 0.275	Depositor DCC
R_{free} test set	1147 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	88.2	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8077	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: 8OG, DDG, MN, DTP, 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	0.41	0/7552	0.55	1/10205 (0.0%)
2	P	0.32	0/297	0.89	2/457 (0.4%)
3	T	1.00	1/383 (0.3%)	1.06	3/586 (0.5%)
All	All	0.45	1/8232 (0.0%)	0.61	6/11248 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	T	1	DC	O3'-P	-18.58	1.38	1.61

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	1	DC	OP1-P-O3'	-11.60	79.69	105.20
3	T	1	DC	P-O3'-C3'	10.60	132.42	119.70
3	T	1	DC	O3'-P-O5'	9.24	121.56	104.00
1	A	170	LEU	CA-CB-CG	5.38	127.67	115.30
2	P	105	DC	P-O3'-C3'	5.37	126.14	119.70
2	P	112	DA	O5'-P-OP2	-5.17	101.05	105.70

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	7371	0	7261	111	0
2	P	286	0	159	2	0
3	T	366	0	203	6	0
4	A	30	0	12	2	0
5	A	3	0	0	0	0
6	P	5	0	0	0	0
6	T	5	0	0	0	0
7	A	10	0	0	0	0
7	P	1	0	0	0	0
All	All	8077	0	7635	117	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 7.

All (117) 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:475:ILE:HG13	1:A:566:LEU:HD23	1.61	0.82
1:A:112:ASN:HB3	1:A:214:THR:HG23	1.63	0.81
1:A:424:ASN:O	1:A:429:THR:HG21	1.81	0.81
1:A:685:ARG:NH1	1:A:686:GLU:O	2.15	0.78
1:A:441:ASP:HB3	1:A:447:ALA:HB2	1.70	0.74
1:A:410:PHE:HB3	1:A:683:MET:HG2	1.70	0.73
1:A:171:GLN:H	1:A:177:GLU:HG3	1.53	0.73
1:A:560:LYS:HA	1:A:563:ILE:HD11	1.71	0.70
1:A:112:ASN:HB3	1:A:214:THR:CG2	2.22	0.70
1:A:839:ASN:HD22	1:A:841:PHE:HD2	1.39	0.69
1:A:415:LEU:HD22	1:A:623:ASP:HB3	1.76	0.68
1:A:546:GLN:O	1:A:550:VAL:HG23	1.94	0.67
1:A:579:ASP:OD1	1:A:581:ARG:HB2	1.95	0.67
1:A:689:ALA:HB2	1:A:712:VAL:HG22	1.76	0.66
1:A:126:PRO:HA	1:A:225:TYR:HD2	1.60	0.66
1:A:129:ALA:HA	1:A:225:TYR:CE2	2.31	0.65
1:A:705:LYS:HD3	3:T:7:DA:H5"	1.80	0.64
1:A:897:LEU:HD22	1:A:897:LEU:H	1.64	0.61
1:A:382:GLN:HG2	1:A:383:GLY:N	2.15	0.60
1:A:15:ILE:HD13	1:A:65:MET:CE	2.32	0.59
1:A:560:LYS:O	1:A:563:ILE:HG12	2.05	0.57
3:T:3:8OG:H2"	3:T:4:DC:C6	2.40	0.57
2:P:113:DA:H2"	2:P:114:DDG:H5"	1.86	0.57
1:A:560:LYS:HA	1:A:563:ILE:CD1	2.35	0.56
1:A:116:GLU:HB2	1:A:135:ALA:HB3	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:LEU:HD11	1:A:470:VAL:HG21	1.87	0.55
1:A:295:GLU:O	1:A:295:GLU:HG3	2.07	0.54
1:A:94:SER:OG	1:A:374:LYS:HD2	2.07	0.54
1:A:391:TYR:HB2	1:A:392:PRO:HD2	1.90	0.54
1:A:417:PRO:HA	1:A:420:ILE:HD12	1.89	0.54
1:A:415:LEU:HD22	1:A:623:ASP:CB	2.37	0.53
1:A:633:ILE:HD11	1:A:651:LEU:HD11	1.89	0.53
1:A:644:THR:O	1:A:648:VAL:HG23	2.09	0.53
1:A:848:TRP:CE2	1:A:854:ILE:HG12	2.44	0.53
1:A:149:PHE:HB3	1:A:197:LEU:CD1	2.38	0.52
1:A:223:ILE:HB	1:A:224:PRO:HD3	1.91	0.52
1:A:412:LEU:HB2	1:A:623:ASP:HB2	1.91	0.52
1:A:523:SER:H	1:A:526:ILE:HD12	1.75	0.52
1:A:424:ASN:O	1:A:429:THR:CG2	2.54	0.52
1:A:428:GLU:OE2	1:A:470:VAL:HG23	2.09	0.52
1:A:839:ASN:ND2	1:A:841:PHE:HB2	2.24	0.52
1:A:304:LYS:O	1:A:323:TYR:OH	2.28	0.51
1:A:458:PRO:HG2	1:A:588:THR:HG22	1.91	0.51
1:A:715:MET:O	1:A:715:MET:HG3	2.10	0.51
1:A:116:GLU:HG2	1:A:324:ASN:ND2	2.25	0.51
1:A:382:GLN:HG2	1:A:383:GLY:H	1.74	0.51
1:A:429:THR:HG23	1:A:463:TYR:HD1	1.76	0.51
1:A:222:ALA:O	1:A:226:VAL:HG12	2.11	0.50
1:A:249:ARG:HB2	1:A:249:ARG:HH11	1.77	0.50
1:A:835:LEU:HD11	1:A:846:ILE:HB	1.95	0.49
1:A:414:SER:HB3	1:A:417:PRO:HG2	1.94	0.49
1:A:126:PRO:HA	1:A:225:TYR:CD2	2.46	0.48
1:A:821:ALA:HB2	1:A:857:LEU:HD12	1.95	0.48
1:A:198:LEU:HD23	1:A:234:PHE:HE2	1.78	0.48
1:A:831:TYR:O	1:A:847:ALA:HA	2.13	0.48
1:A:440:HIS:HA	1:A:443:ILE:HD12	1.96	0.48
1:A:417:PRO:HB3	1:A:475:ILE:HD13	1.96	0.48
1:A:482:ARG:NH1	4:A:904:DTP:O1G	2.45	0.47
3:T:8:DT:H2'	3:T:9:DG:C8	2.49	0.47
1:A:373:LEU:HD11	1:A:470:VAL:CG2	2.44	0.47
1:A:499:ILE:O	1:A:503:LEU:HD23	2.15	0.47
1:A:685:ARG:NH2	1:A:717:GLY:H	2.12	0.47
1:A:171:GLN:OE1	1:A:319:ARG:NH2	2.48	0.47
1:A:688:ILE:HD12	1:A:714:ASP:HB3	1.96	0.47
1:A:298:LEU:HD11	1:A:334:ILE:HG13	1.97	0.46
1:A:872:LEU:HD12	1:A:876:PHE:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:PHE:HB3	1:A:197:LEU:HD12	1.96	0.46
1:A:163:SER:OG	1:A:166:ILE:HG12	2.17	0.46
1:A:412:LEU:HD23	1:A:415:LEU:HD13	1.96	0.46
1:A:544:ARG:O	1:A:548:THR:OG1	2.35	0.45
1:A:725:LEU:HD11	1:A:750:ARG:HB2	1.97	0.45
1:A:831:TYR:CD2	1:A:850:SER:HA	2.52	0.45
4:A:904:DTP:H8	2:P:114:DDG:H2"	1.99	0.45
1:A:113:PHE:HA	1:A:137:THR:O	2.16	0.45
1:A:33:TYR:HB3	1:A:65:MET:CE	2.47	0.45
1:A:818:ASN:H	1:A:818:ASN:HD22	1.65	0.45
1:A:401:PRO:HA	1:A:702:TRP:O	2.17	0.44
3:T:3:8OG:H2"	3:T:4:DC:H6	1.81	0.44
1:A:326:ILE:O	1:A:330:ARG:HG2	2.18	0.44
1:A:457:SER:OG	1:A:458:PRO:HD2	2.17	0.44
1:A:230:ILE:HG23	1:A:234:PHE:HD2	1.83	0.44
1:A:706:LYS:O	1:A:729:GLY:HA3	2.18	0.44
1:A:171:GLN:N	1:A:177:GLU:HG3	2.26	0.44
1:A:655:ALA:HA	1:A:659:MET:HB2	1.99	0.44
1:A:397:LYS:O	1:A:399:PRO:HD3	2.18	0.44
1:A:702:TRP:CZ3	1:A:710:LEU:HD21	2.52	0.44
1:A:707:ARG:HD2	3:T:8:DT:H4'	1.99	0.43
1:A:449:ARG:HA	1:A:450:PRO:HD3	1.90	0.43
1:A:702:TRP:CH2	1:A:710:LEU:HD21	2.53	0.43
1:A:422:GLN:O	1:A:676:ASN:HB3	2.19	0.43
1:A:415:LEU:CD2	1:A:623:ASP:HB3	2.47	0.43
1:A:641:PHE:CD1	1:A:647:TRP:HB3	2.54	0.43
1:A:420:ILE:HG21	1:A:471:VAL:HG12	2.01	0.43
1:A:422:GLN:NE2	1:A:671:CYS:SG	2.92	0.43
1:A:51:ASP:OD1	1:A:51:ASP:C	2.57	0.42
1:A:494:ARG:HD2	1:A:521:ASP:OD1	2.18	0.42
1:A:202:LEU:HD23	1:A:202:LEU:HA	1.74	0.42
1:A:426:SER:OG	1:A:427:PRO:HD2	2.19	0.42
1:A:593:ALA:HA	1:A:670:MET:CE	2.49	0.42
3:T:6:DT:C2	3:T:7:DA:C8	3.07	0.42
1:A:15:ILE:HD13	1:A:65:MET:HE3	1.99	0.42
1:A:707:ARG:NH2	1:A:731:GLU:OE1	2.50	0.42
1:A:76:GLU:HG2	1:A:382:GLN:HE22	1.84	0.42
1:A:730:LEU:HD22	1:A:883:PHE:CE1	2.54	0.42
1:A:415:LEU:O	1:A:419:ILE:HG13	2.19	0.42
1:A:481:GLN:HB3	1:A:559:ARG:HD3	2.00	0.42
1:A:737:THR:HB	1:A:742:GLN:HE21	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:248:THR:HA	1:A:264:THR:O	2.20	0.41
1:A:856:ASP:C	1:A:858:ILE:H	2.23	0.41
1:A:229:ARG:HD2	1:A:229:ARG:HA	1.92	0.41
1:A:635:LYS:HD2	1:A:635:LYS:HA	1.96	0.41
1:A:751:ARG:NH1	1:A:763:TYR:HB2	2.36	0.41
1:A:230:ILE:HG23	1:A:234:PHE:CD2	2.56	0.41
1:A:7:THR:HG21	1:A:267:GLY:O	2.20	0.41
1:A:85:MET:HB2	1:A:370:PHE:CE1	2.56	0.41
1:A:647:TRP:O	1:A:650:PHE:HB3	2.22	0.40
1:A:337:LYS:HG3	1:A:547:ARG:HD2	2.03	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	901/903 (100%)	837 (93%)	58 (6%)	6 (1%)	25 67

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	820	ASP
1	A	858	ILE
1	A	424	ASN
1	A	303	LEU
1	A	716	GLU
1	A	42	PRO

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	799/799 (100%)	752 (94%)	47 (6%)	23	60

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	LYS
1	A	14	SER
1	A	15	ILE
1	A	52	ILE
1	A	58	THR
1	A	77	ASP
1	A	107	LYS
1	A	121	ASP
1	A	160	GLU
1	A	189	MET
1	A	220	SER
1	A	249	ARG
1	A	276	LEU
1	A	284	ASN
1	A	295	GLU
1	A	302	LYS
1	A	373	LEU
1	A	389	GLN
1	A	412	LEU
1	A	429	THR
1	A	468	ASP
1	A	475	ILE
1	A	505	ASN
1	A	510	VAL
1	A	517	ASP
1	A	525	GLU
1	A	530	ILE
1	A	548	THR
1	A	563	ILE

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Mol	Chain	Res	Type
1	A	567	TYR
1	A	580	LEU
1	A	581	ARG
1	A	607	GLU
1	A	618	LEU
1	A	635	LYS
1	A	639	SER
1	A	641	PHE
1	A	658	ARG
1	A	739	LYS
1	A	743	LYS
1	A	818	ASN
1	A	843	ASP
1	A	854	ILE
1	A	855	THR
1	A	897	LEU
1	A	898	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	ASN
1	A	284	ASN
1	A	382	GLN
1	A	422	GLN
1	A	444	ASN
1	A	558	ASN
1	A	646	HIS
1	A	676	ASN
1	A	786	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
2	DDG	P	114	3,2	16,23,24	1.37	3 (18%)	15,33,36	3.01	9 (60%)
3	8OG	T	3	3	17,25,26	1.43	2 (11%)	22,37,40	4.07	7 (31%)

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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DDG	P	114	3,2	-	0/3/18/19	0/3/3/3
3	8OG	T	3	3	-	0/3/21/22	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	114	DDG	O5'-C5'	-2.02	1.41	1.44
2	P	114	DDG	C5-C4	2.81	1.46	1.40
2	P	114	DDG	C6-C5	3.79	1.48	1.41
3	T	3	8OG	C5-C4	3.88	1.49	1.40
3	T	3	8OG	C6-C5	3.89	1.48	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	3	8OG	C2'-C1'-N9	-13.85	101.73	115.79
3	T	3	8OG	C6-C5-C4	-4.54	116.33	120.84
2	P	114	DDG	C5-C6-N1	-4.32	117.33	123.48
2	P	114	DDG	C2'-C1'-N9	-4.18	104.50	112.47
3	T	3	8OG	C5-C6-N1	-3.94	117.88	123.48
2	P	114	DDG	C4-C5-N7	-3.73	105.80	109.41
2	P	114	DDG	N3-C2-N1	-3.39	122.51	127.46
3	T	3	8OG	N3-C2-N1	-3.31	122.62	127.46
2	P	114	DDG	C6-C5-C4	-3.22	117.64	120.84
2	P	114	DDG	O4'-C1'-N9	2.31	111.68	107.78
2	P	114	DDG	C3'-C2'-C1'	2.41	105.36	102.69
3	T	3	8OG	C6-N1-C2	4.75	122.89	116.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	114	DDG	C6-N1-C2	4.75	122.89	116.06
2	P	114	DDG	C2-N3-C4	4.83	120.80	115.16
3	T	3	8OG	C2-N3-C4	4.94	120.92	115.16
3	T	3	8OG	O4'-C1'-N9	8.09	115.70	108.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	114	DDG	2	0
3	T	3	8OG	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	DTP	A	904	5	26,32,32	1.00	1 (3%)	26,50,50	1.91	4 (15%)
6	SO4	P	2	-	4,4,4	0.43	0	6,6,6	0.19	0
6	SO4	T	19	-	4,4,4	0.44	0	6,6,6	0.24	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
4	DTP	A	904	5	-	0/18/34/34	0/3/3/3
6	SO4	P	2	-	-	0/0/0/0	0/0/0/0
6	SO4	T	19	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	904	DTP	C5-C4	3.35	1.48	1.40

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	904	DTP	N3-C2-N1	-6.52	123.18	128.86
4	A	904	DTP	C2'-C1'-N9	-4.21	104.28	114.23
4	A	904	DTP	O4'-C1'-N9	2.14	111.39	107.78
4	A	904	DTP	N6-C6-N1	2.14	123.02	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	904	DTP	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	T	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	T	1:DC	O3'	2:DA	P	1.38

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	A	903/903 (100%)	-0.24	13 (1%) 75 49	52, 91, 151, 218	0
2	P	13/14 (92%)	-0.01	1 (7%) 14 5	53, 68, 133, 134	0
3	T	16/18 (88%)	-0.44	0 100 100	51, 61, 97, 100	0
All	All	932/935 (99%)	-0.24	14 (1%) 74 47	51, 91, 151, 218	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	44	SER	4.9
1	A	32	GLU	3.9
1	A	254	GLU	3.7
1	A	43	GLU	2.9
1	A	517	ASP	2.7
2	P	101	DG	2.5
1	A	259	SER	2.4
1	A	523	SER	2.3
1	A	256	MET	2.3
1	A	31	VAL	2.2
1	A	516	VAL	2.2
1	A	258	GLY	2.2
1	A	253	ILE	2.0
1	A	255	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	DDG	P	114	21/22	0.98	0.17	-	43,54,58,59	0
3	8OG	T	3	23/24	0.97	0.11	-	42,55,62,66	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
6	SO4	P	2	5/5	0.85	0.20	-0.21	107,109,118,120	0
6	SO4	T	19	5/5	0.94	0.14	-0.66	71,75,77,84	0
4	DTP	A	904	30/30	0.97	0.16	-1.18	40,46,55,56	0
5	MN	A	908	1/1	0.94	0.13	-3.13	76,76,76,76	0
5	MN	A	907	1/1	0.99	0.11	-6.69	53,53,53,53	0
5	MN	A	906	1/1	0.98	0.18	-	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.