



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 16, 2023 – 05:23 PM EST

PDB ID : 4PEI  
Title : Dbr1 in complex with synthetic branched RNA analog  
Authors : Montemayor, E.J.; Katolik, A.; Clark, N.E.; Taylor, A.B.; Schuermann, J.P.; Combs, D.J.; Johnsson, R.; Holloway, S.P.; Stevens, S.W.; Damha, M.J.; Hart, P.J.  
Deposited on : 2014-04-23  
Resolution : 1.95 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

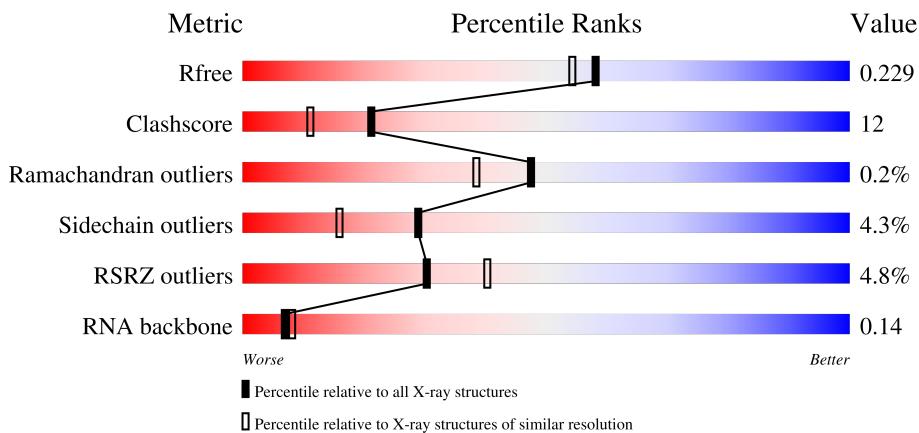
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

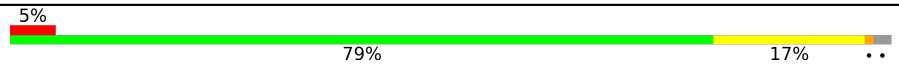
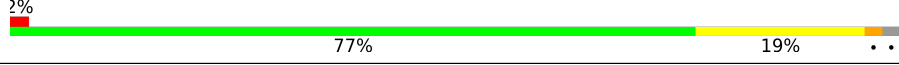
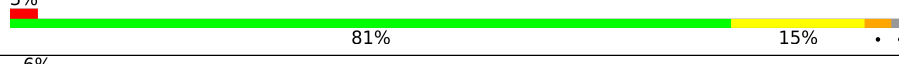

The reported resolution of this entry is 1.95 Å.

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))
$R_{free}$	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)
RNA backbone	3102	1124 (2.50-1.42)

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  $\geq 3$ , 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	356	 5% 79% 17% ..
1	B	356	 2% 77% 19% ..
1	C	356	 3% 81% 15% ..
1	D	356	 6% 79% 16% ..

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Mol	Chain	Length	Quality of chain
1	E	356	
2	V	5	
2	W	5	
2	X	5	
2	Y	5	
2	Z	5	
3	Q	2	
3	R	2	
3	S	2	
3	T	2	
3	U	2	

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
3	G46	Q	510	-	-	X	-
3	G46	R	510	-	-	X	-
3	G46	T	510	-	-	X	-
3	G46	U	501	-	-	-	X
7	PG4	B	403	-	-	X	-

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 16230 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 RNA lariat debranching enzyme, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	349	2872	1866	467	525	14	0	3	0
1	B	349	2876	1869	469	524	14	0	3	0
1	C	349	2869	1864	466	525	14	0	2	0
1	D	349	2873	1866	469	524	14	0	2	0
1	E	348	2875	1867	470	524	14	0	3	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP C4M1P9
A	0	ALA	-	expression tag	UNP C4M1P9
A	14	SER	CYS	engineered mutation	UNP C4M1P9
B	-1	GLY	-	expression tag	UNP C4M1P9
B	0	ALA	-	expression tag	UNP C4M1P9
B	14	SER	CYS	engineered mutation	UNP C4M1P9
C	-1	GLY	-	expression tag	UNP C4M1P9
C	0	ALA	-	expression tag	UNP C4M1P9
C	14	SER	CYS	engineered mutation	UNP C4M1P9
D	-1	GLY	-	expression tag	UNP C4M1P9
D	0	ALA	-	expression tag	UNP C4M1P9
D	14	SER	CYS	engineered mutation	UNP C4M1P9
E	-1	GLY	-	expression tag	UNP C4M1P9
E	0	ALA	-	expression tag	UNP C4M1P9
E	14	SER	CYS	engineered mutation	UNP C4M1P9

- Molecule 2 is a RNA chain called RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3').

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	V	3	Total	C	N	O	P	0	0	0
			32	14	5	11	2			
2	W	3	Total	C	N	O	P	0	0	0
			32	14	5	11	2			
2	X	3	Total	C	N	O	P	0	0	0
			63	29	13	18	3			
2	Y	2	Total	C	N	O	P	0	0	1
			22	10	5	6	1			
2	Z	3	Total	C	N	O	P	0	0	0
			58	28	13	15	2			

- Molecule 3 is a RNA chain called RNA (5'-R\*(G46)P\*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
3	Q	1	Total	C	N	O	P	S	0	0	0
			24	10	5	7	1	1			
3	R	1	Total	C	O	P	S		0	0	0
			13	5	6	1	1				
3	S	1	Total	O	P	S			0	0	0
			5	3	1	1					
3	T	1	Total	O	P	S			0	0	0
			5	3	1	1					
3	U	1	Total	O					0	0	1
			1	1							

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

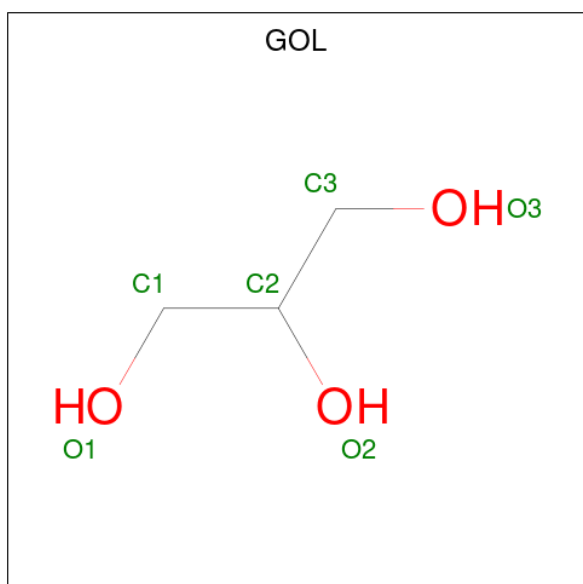
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ni	0	0
			1	1		
4	B	1	Total	Ni	0	0
			1	1		
4	C	1	Total	Ni	0	0
			1	1		
4	D	1	Total	Ni	0	0
			1	1		
4	E	1	Total	Ni	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



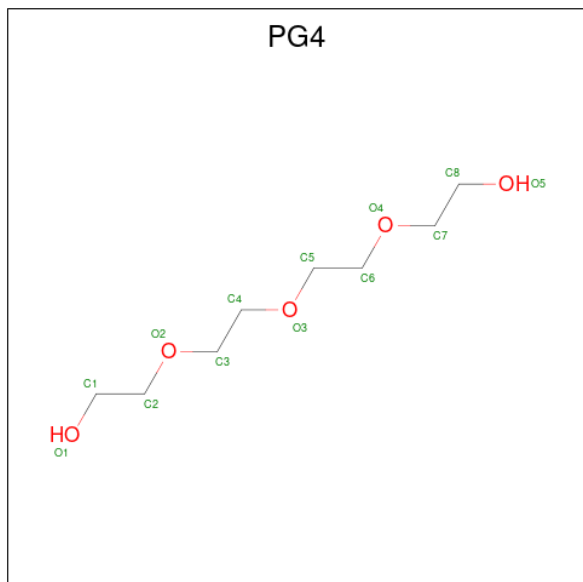
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O S 5 4 1	0	0
5	B	1	Total O S 5 4 1	0	0
5	C	1	Total O S 5 4 1	0	0
5	D	1	Total O S 5 4 1	0	0
5	E	1	Total O S 5 4 1	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	338	Total	O	0	0
			338	338		
8	V	3	Total	O	0	0
			3	3		
8	Q	2	Total	O	0	0
			2	2		

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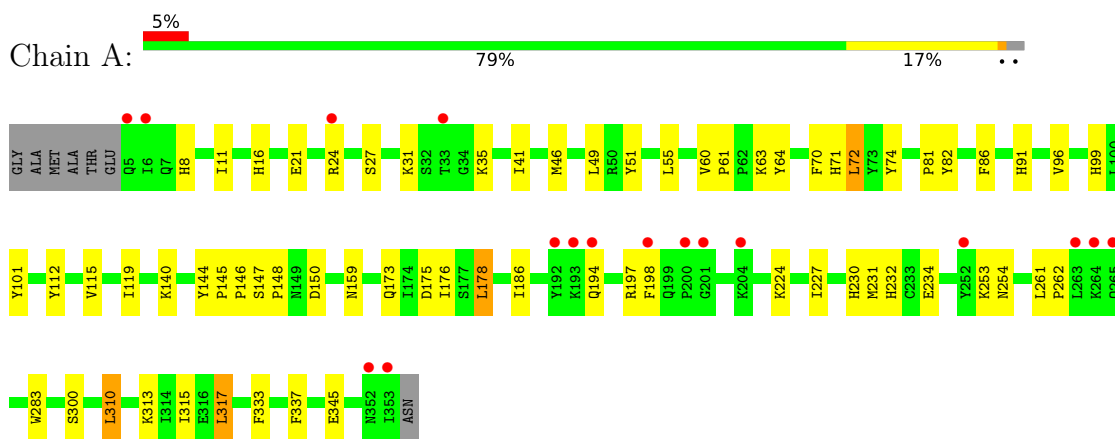
<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
8	B	271	Total O 271 271	0	0
8	W	2	Total O 2 2	0	0
8	R	1	Total O 1 1	0	0
8	C	325	Total O 325 325	0	0
8	X	5	Total O 5 5	0	0
8	D	301	Total O 301 301	0	0
8	Y	4	Total O 4 4	0	0
8	E	277	Total O 277 277	0	0
8	Z	2	Total O 2 2	0	0



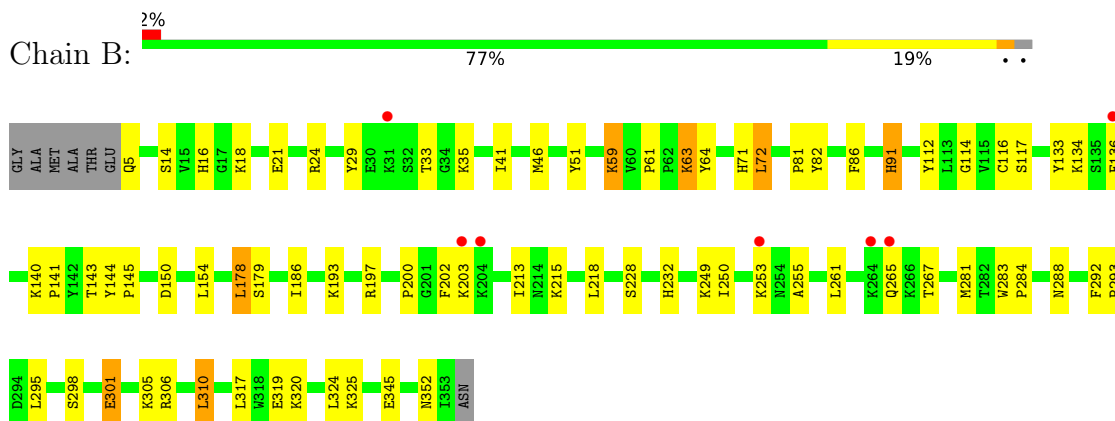
### 3 Residue-property plots [i](#)

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.

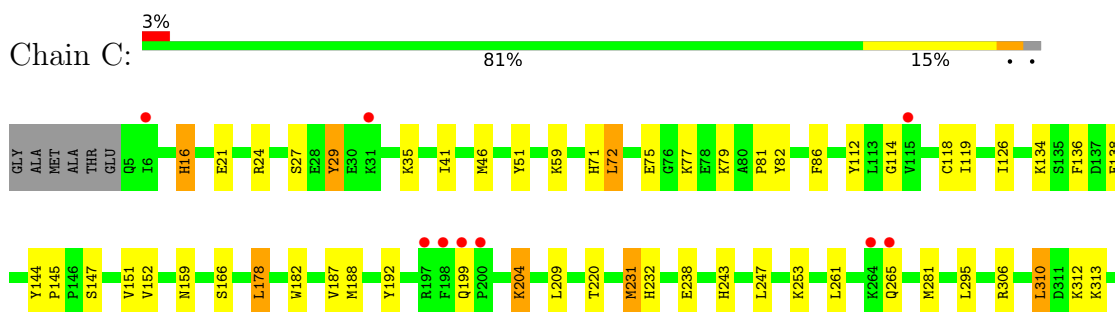
- Molecule 1: RNA lariat debranching enzyme, putative



- Molecule 1: RNA lariat debranching enzyme, putative

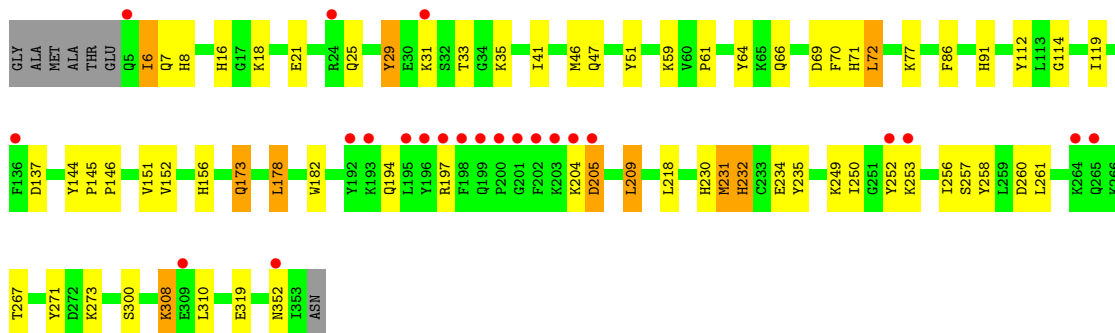
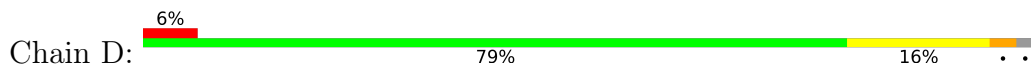


- Molecule 1: RNA lariat debranching enzyme, putative

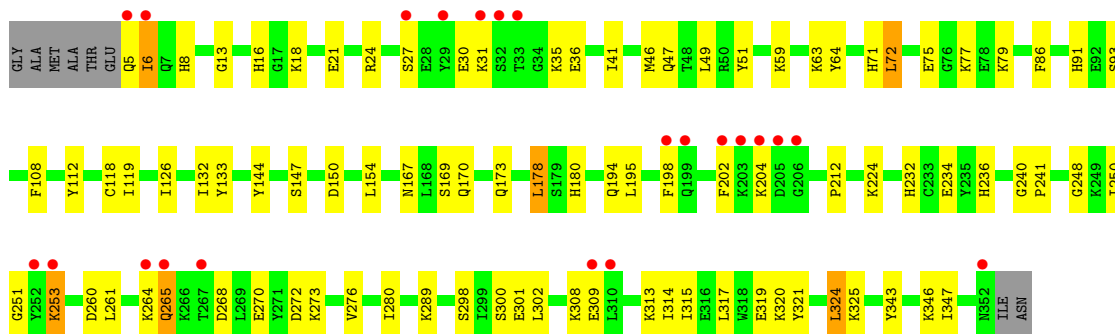




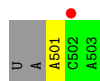
- Molecule 1: RNA lariat debranching enzyme, putative



- Molecule 1: RNA lariat debranching enzyme, putative




- Molecule 2: RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3')



- Molecule 2: RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3')



- Molecule 2: RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3')

Chain X:  20% 40% 40%



• Molecule 2: RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3')

Chain Y:  40% 20% 60%



• Molecule 2: RNA (5'-R(\*UP\*AP\*AP\*CP\*A)-3')

Chain Z:  20% 60% 40%



• Molecule 3: RNA (5'-R(\*(G46)P\*U)-3')

Chain Q:  50% 50%



• Molecule 3: RNA (5'-R(\*(G46)P\*U)-3')

Chain R:  50% 50%



• Molecule 3: RNA (5'-R(\*(G46)P\*U)-3')

Chain S:  50% 50%



• Molecule 3: RNA (5'-R(\*(G46)P\*U)-3')

Chain T:  50% 50%



• Molecule 3: RNA (5'-R(\*(G46)P\*U)-3')

Chain U:  50% 50%

CSG1  
0

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.04Å 142.48Å 214.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.37 – 1.95 46.37 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.8 (46.37-1.95) 98.8 (46.37-1.95)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 1.95Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.2_1309)	Depositor
R, $R_{free}$	0.184 , 0.229 0.184 , 0.229	Depositor DCC
$R_{free}$ test set	8110 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.5	Xtrriage
Anisotropy	0.247	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 51.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	16230	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: SO4, GOL, NI, G46, 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	A	0.68	0/2956	0.71	0/3995
1	B	0.69	1/2961 (0.0%)	0.72	0/4002
1	C	0.72	0/2950	0.73	0/3987
1	D	0.66	0/2952	0.71	0/3990
1	E	0.68	1/2958 (0.0%)	0.72	1/3998 (0.0%)
2	V	0.65	0/35	1.01	0/53
2	W	0.90	0/35	1.15	0/53
2	X	0.63	0/70	1.13	1/106 (0.9%)
2	Y	2.37	1/24 (4.2%)	2.09	2/36 (5.6%)
2	Z	0.53	0/65	0.91	0/99
All	All	0.69	3/15006 (0.0%)	0.73	4/20319 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Y	500	A	O3'-P	-10.65	1.48	1.61
1	E	241	PRO	N-CD	5.43	1.55	1.47
1	B	91	HIS	C-N	-5.10	1.22	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Y	501	A	OP1-P-OP2	-8.10	107.45	119.60
2	X	501	A	C4'-C3'-C2'	-5.96	96.64	102.60
2	Y	500	A	OP1-P-O3'	5.57	117.45	105.20
1	E	240	GLY	C-N-CD	5.53	140.01	128.40

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	2872	0	2846	58	0
1	B	2876	0	2851	73	0
1	C	2869	0	2842	54	0
1	D	2873	0	2842	73	0
1	E	2875	0	2839	83	0
2	V	32	0	14	2	0
2	W	32	0	14	4	0
2	X	63	0	34	3	0
2	Y	22	0	12	17	0
2	Z	58	0	30	13	0
3	Q	24	0	13	7	0
3	R	13	0	8	7	0
3	S	5	0	1	0	0
3	T	5	0	1	7	0
3	U	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	1	0
5	E	5	0	0	0	0
6	A	12	0	16	2	0
6	C	12	0	16	2	0
6	D	12	0	16	1	0
7	B	13	0	18	14	0
8	A	338	0	0	12	0
8	B	271	0	0	9	0
8	C	325	0	0	8	0
8	D	301	0	0	11	0
8	E	277	0	0	7	0
8	Q	2	0	0	0	0
8	R	1	0	0	1	0
8	V	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	W	2	0	0	0	0
8	X	5	0	0	0	0
8	Y	4	0	0	1	0
8	Z	2	0	0	0	0
All	All	16230	0	14413	352	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 12.

The worst 5 of 352 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:E:91[B]:HIS:NE2	2:Z:502:C:O2'	1.57	1.38
1:D:91:HIS:CE1	2:Y:501:A:H1'	1.62	1.34
1:E:91[B]:HIS:CE1	2:Z:502:C:O2'	1.91	1.24
1:B:143:THR:HG23	7:B:403:PG4:H41	1.24	1.17
1:B:16[B]:HIS:NE2	3:R:510:G46:S2P	2.20	1.13

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	350/356 (98%)	342 (98%)	7 (2%)	1 (0%)	41	30
1	B	350/356 (98%)	342 (98%)	7 (2%)	1 (0%)	41	30
1	C	349/356 (98%)	341 (98%)	8 (2%)	0	100	100
1	D	349/356 (98%)	337 (97%)	11 (3%)	1 (0%)	41	30
1	E	349/356 (98%)	338 (97%)	11 (3%)	0	100	100
All	All	1747/1780 (98%)	1700 (97%)	44 (2%)	3 (0%)	47	38



All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	230	HIS
1	D	232	HIS
1	B	352	ASN

### 5.3.2 Protein sidechains [i](#)

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	318/320 (99%)	310 (98%)	8 (2%)	47	38
1	B	319/320 (100%)	306 (96%)	13 (4%)	30	18
1	C	318/320 (99%)	302 (95%)	16 (5%)	24	11
1	D	318/320 (99%)	301 (95%)	17 (5%)	22	10
1	E	318/320 (99%)	304 (96%)	14 (4%)	28	15
All	All	1591/1600 (99%)	1523 (96%)	68 (4%)	29	16

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	72	LEU
1	E	169	SER
1	E	300	SER
1	C	51	TYR
1	C	29	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	265	GLN
1	E	163	GLN
1	D	173	GLN
1	C	232	HIS
1	E	47	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	V	0/5	-	-
2	W	1/5 (20%)	0	0
2	X	2/5 (40%)	1 (50%)	0
2	Y	0/5	-	-
2	Z	1/5 (20%)	0	0
3	Q	0/2	-	-
3	R	0/2	-	-
3	S	0/2	-	-
3	T	0/2	-	-
3	U	0/2	-	-
All	All	4/35 (11%)	1 (25%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	X	502	C

There are no RNA pucker outliers to report.

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

Of 5 non-standard protein/DNA/RNA residues modelled in this entry, 1 is modelled with single atom - 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
3	G46	Q	510	2,4	20,26,26	1.39	4 (20%)	22,40,40	1.29	2 (9%)
3	G46	R	510	2,4	11,13,26	1.25	1 (9%)	13,19,40	1.42	2 (15%)
3	G46	T	510	2,4	2,4,26	2.39	1 (50%)	3,6,40	3.11	1 (33%)
3	G46	S	510	2	2,4,26	2.72	2 (100%)	3,6,40	2.70	1 (33%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	G46	Q	510	2,4	-	1/5/26/26	0/3/3/3
3	G46	R	510	2,4	-	1/5/19/26	0/1/1/3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	R	510	G46	P-O1P	-3.25	1.48	1.56
3	S	510	G46	P-O1P	-3.24	1.48	1.56
3	Q	510	G46	P-O1P	-3.20	1.48	1.56
3	T	510	G46	P-O1P	-3.18	1.48	1.56
3	Q	510	G46	C6-N1	-2.26	1.34	1.37

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	510	G46	O5'-P-O3P	-5.08	100.20	113.38
3	S	510	G46	O5'-P-O3P	-4.17	102.57	113.38
3	Q	510	G46	C3'-C2'-C1'	3.42	106.13	100.98
3	R	510	G46	O4'-C4'-C3'	3.02	107.38	104.70
3	Q	510	G46	C8-N7-C5	2.41	107.57	102.99

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Q	510	G46	O4'-C4'-C5'-O5'
3	R	510	G46	O4'-C4'-C5'-O5'

There are no ring outliers.

3 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	Q	510	G46	7	0
3	R	510	G46	7	0
3	T	510	G46	7	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 5 are monoatomic - leaving 12 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
5	SO4	D	402	-	4,4,4	0.22	0	6,6,6	0.72	0
5	SO4	B	402	-	4,4,4	0.12	0	6,6,6	0.40	0
6	GOL	A	403	-	5,5,5	0.31	0	5,5,5	0.48	0
6	GOL	C	403	-	5,5,5	0.34	0	5,5,5	0.69	0
5	SO4	A	402	-	4,4,4	0.82	0	6,6,6	1.38	1 (16%)
6	GOL	C	404	-	5,5,5	0.20	0	5,5,5	1.73	2 (40%)
6	GOL	A	404	-	5,5,5	0.53	0	5,5,5	0.39	0
6	GOL	D	404	-	5,5,5	0.33	0	5,5,5	0.75	0
5	SO4	C	402	-	4,4,4	0.18	0	6,6,6	0.51	0
5	SO4	E	402	-	4,4,4	0.16	0	6,6,6	0.23	0
7	PG4	B	403	-	12,12,12	0.64	0	11,11,11	0.39	0
6	GOL	D	403	-	5,5,5	0.32	0	5,5,5	0.66	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
6	GOL	A	403	-	-	1/4/4/4	-
6	GOL	C	403	-	-	3/4/4/4	-
6	GOL	C	404	-	-	2/4/4/4	-
6	GOL	A	404	-	-	0/4/4/4	-
6	GOL	D	404	-	-	4/4/4/4	-
7	PG4	B	403	-	-	4/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	D	403	-	-	0/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	402	SO4	O3-S-O1	3.08	125.40	109.31
6	C	404	GOL	C3-C2-C1	-2.82	100.76	111.70
6	C	404	GOL	O1-C1-C2	-2.46	98.42	110.20

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	403	GOL	C1-C2-C3-O3
6	D	404	GOL	O1-C1-C2-C3
6	D	404	GOL	C1-C2-C3-O3
7	B	403	PG4	O1-C1-C2-O2
7	B	403	PG4	O2-C3-C4-O3

There are no ring outliers.

5 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	402	SO4	1	0
6	C	404	GOL	2	0
6	A	404	GOL	2	0
6	D	404	GOL	1	0
7	B	403	PG4	14	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	349/356 (98%)	0.21	17 (4%) 29 39	13, 22, 39, 55	0
1	B	349/356 (98%)	0.09	7 (2%) 65 73	13, 22, 39, 57	0
1	C	349/356 (98%)	0.25	11 (3%) 47 57	11, 21, 36, 57	0
1	D	349/356 (98%)	0.31	23 (6%) 18 26	11, 22, 45, 63	0
1	E	348/356 (97%)	0.35	22 (6%) 20 28	12, 25, 44, 58	0
2	V	3/5 (60%)	2.00	1 (33%) 0 0	39, 39, 40, 44	3 (100%)
2	W	3/5 (60%)	2.08	1 (33%) 0 0	37, 37, 42, 43	3 (100%)
2	X	3/5 (60%)	0.64	0 100 100	30, 30, 32, 45	2 (66%)
2	Y	2/5 (40%)	2.60	2 (100%) 0 0	47, 47, 47, 59	2 (100%)
2	Z	3/5 (60%)	1.66	1 (33%) 0 0	35, 35, 40, 49	3 (100%)
3	Q	0/2	-	-	-	-
3	R	0/2	-	-	-	-
3	S	0/2	-	-	-	-
3	T	0/2	-	-	-	-
3	U	0/2	-	-	-	-
All	All	1758/1815 (96%)	0.25	85 (4%) 30 40	11, 22, 42, 63	13 (0%)

The worst 5 of 85 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	198	PHE	7.5
1	E	198	PHE	7.5
1	A	198	PHE	5.5
1	D	202	PHE	4.8
1	E	264	LYS	4.6

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	G46	U	501	1/24	0.27	0.44	48,48,48,48	1
3	G46	S	510	5/24	0.60	0.25	47,50,58,62	5
3	G46	R	510	13/24	0.82	0.21	19,38,45,46	13
3	G46	Q	510	24/24	0.84	0.19	26,45,47,49	24
3	G46	T	510	5/24	0.91	0.13	24,44,46,52	5

## 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
6	GOL	A	404	6/6	0.75	0.25	30,34,43,44	0
6	GOL	D	404	6/6	0.83	0.24	23,32,38,38	0
7	PG4	B	403	13/13	0.84	0.40	24,34,47,50	0
6	GOL	A	403	6/6	0.86	0.17	26,35,37,44	0
6	GOL	C	404	6/6	0.86	0.29	31,36,40,42	0
6	GOL	C	403	6/6	0.87	0.20	32,35,44,45	0
6	GOL	D	403	6/6	0.87	0.19	28,32,44,47	0
5	SO4	B	402	5/5	0.97	0.10	34,35,39,49	0
5	SO4	C	402	5/5	0.97	0.12	33,33,37,43	0
5	SO4	A	402	5/5	0.97	0.10	31,35,37,38	0
5	SO4	E	402	5/5	0.98	0.10	28,29,35,46	0
5	SO4	D	402	5/5	0.98	0.12	28,28,37,47	0
4	NI	A	401	1/1	0.99	0.04	30,30,30,30	0
4	NI	B	401	1/1	0.99	0.05	23,23,23,23	0
4	NI	D	401	1/1	0.99	0.06	26,26,26,26	0
4	NI	E	401	1/1	0.99	0.09	21,21,21,21	0
4	NI	C	401	1/1	1.00	0.06	18,18,18,18	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.