Full wwPDB X-ray Structure Validation Report (i)

May 18, 2020 – 12:40 am BST

PDB ID	:	6H9H
Title	:	Csf5, CRISPR-Cas type IV Cas6 crRNA endonuclease
Authors	:	Pausch, P.; Bange, G.
Deposited on	:	2018-08-04
Resolution	:	1.75 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

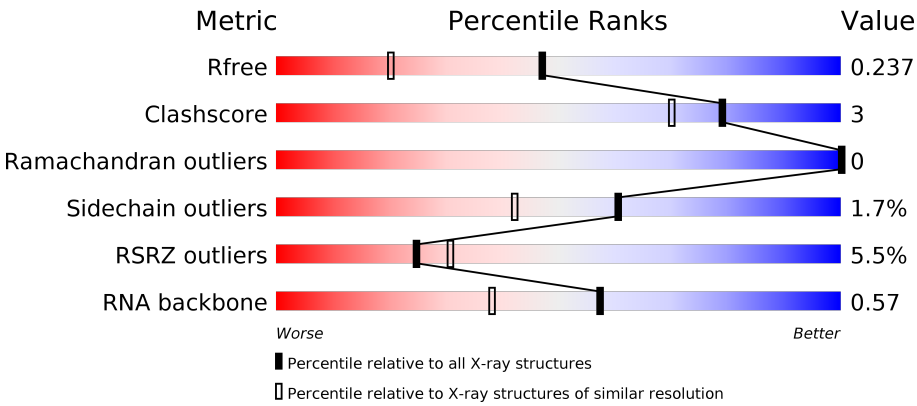
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)
RNA backbone	3102	1079 (2.40-1.12)

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 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 <=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	266	5% 88% 6% 6%
1	B	266	6% 88% 5% 6%
2	D	26	62% 19% 8% 12%
2	E	26	8% 77% 12% • 8%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5345 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 Csf5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	250	1958	1238	363	352	5	0	0	0
1	A	250	1969	1244	367	353	5	0	1	0

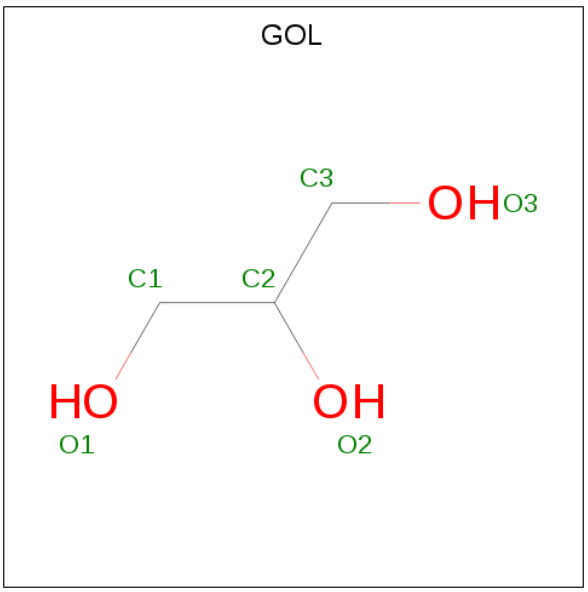
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	initiating methionine	UNP Q5NWP0
B	2	GLY	-	expression tag	UNP Q5NWP0
B	261	HIS	-	expression tag	UNP Q5NWP0
B	262	HIS	-	expression tag	UNP Q5NWP0
B	263	HIS	-	expression tag	UNP Q5NWP0
B	264	HIS	-	expression tag	UNP Q5NWP0
B	265	HIS	-	expression tag	UNP Q5NWP0
B	266	HIS	-	expression tag	UNP Q5NWP0
A	1	MET	-	initiating methionine	UNP Q5NWP0
A	2	GLY	-	expression tag	UNP Q5NWP0
A	261	HIS	-	expression tag	UNP Q5NWP0
A	262	HIS	-	expression tag	UNP Q5NWP0
A	263	HIS	-	expression tag	UNP Q5NWP0
A	264	HIS	-	expression tag	UNP Q5NWP0
A	265	HIS	-	expression tag	UNP Q5NWP0
A	266	HIS	-	expression tag	UNP Q5NWP0

- Molecule 2 is a RNA chain called crRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	D	23	490	217	85	165	23	0	0	0
2	E	24	574	254	98	194	28	0	3	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

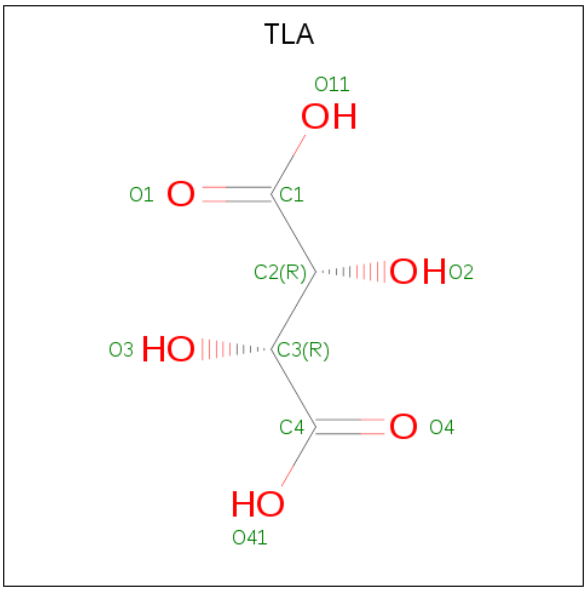


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	B	1	6	3	3	0	0

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mn		
4	B	1	1	1	0	0

- Molecule 5 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	10	4	6	0	0
			Total	C	O		
5	A	1	10	4	6	0	0
			Total	C	O		
5	D	1	10	4	6	0	0
			Total	C	O		
5	E	1	10	4	6	0	0

- Molecule 6 is water.

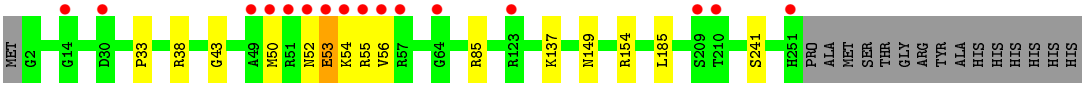
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	B	111	111	111	0	0
			Total	O		
6	A	131	131	131	0	0
			Total	O		
6	D	33	33	33	0	0
			Total	O		
6	E	32	32	32	0	0

3 Residue-property plots (i)

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.

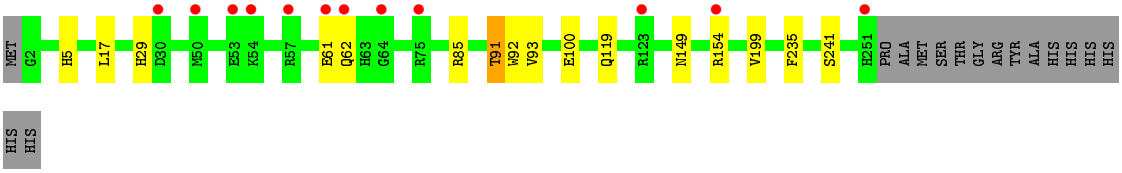
- Molecule 1: Csf5

Chain B:



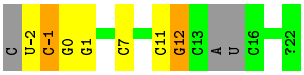
- Molecule 1: Csf5

Chain A:



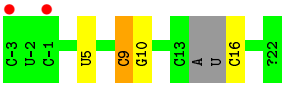
- Molecule 2: crRNA

Chain D:



- Molecule 2: crRNA

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	85.56Å 100.99Å 101.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.25 – 1.75 45.50 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.25-1.75) 93.9 (45.50-1.75)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.94 (at 1.75Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.212 , 0.237 0.212 , 0.237	Depositor DCC
R_{free} test set	4426 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	31.7	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 40.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5345	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL, TLA, MN, 23G

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	0.39	0/2013	0.58	0/2732
1	B	0.39	0/2002	0.59	0/2718
2	D	0.59	0/515	1.28	7/799 (0.9%)
2	E	0.53	0/608	1.17	3/943 (0.3%)
All	All	0.43	0/5138	0.79	10/7192 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	-1	C	N3-C2-O2	-8.44	116.00	121.90
2	D	-1	C	C6-N1-C2	-6.73	117.61	120.30
2	D	-1	C	C5-C4-N4	6.45	124.72	120.20
2	D	-2	U	N3-C2-O2	-6.35	117.76	122.20
2	D	-2	U	C5-C4-O4	6.09	129.56	125.90
2	E	9	C	C6-N1-C2	-5.45	118.12	120.30
2	D	1	G	O4'-C1'-N9	5.33	112.47	108.20
2	E	9	C	C5-C6-N1	5.29	123.64	121.00
2	D	-1	C	N1-C2-N3	5.04	122.73	119.20
2	E	16	C	O4'-C1'-N1	-5.03	104.18	108.20

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	1969	0	1997	12	0
1	B	1958	0	1985	10	0
2	D	490	0	250	4	0
2	E	574	0	292	5	0
3	B	6	0	8	0	0
4	B	1	0	0	0	0
5	A	20	0	8	3	0
5	D	10	0	4	0	0
5	E	10	0	4	0	0
6	A	131	0	0	3	0
6	B	111	0	0	0	0
6	D	33	0	0	0	0
6	E	32	0	0	0	0
All	All	5345	0	4548	27	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 3.

All (27) 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:B:55:ARG:HH22	2:E:10:G:H4'	1.28	0.94
2:D:11:C:H2'	2:D:12:G:H5"	1.59	0.84
1:A:85:ARG:NH2	5:A:301:TLA:O1	2.21	0.74
1:B:52:ASN:HD21	1:B:54:LYS:HB3	1.57	0.70
1:A:85:ARG:HH21	5:A:301:TLA:C1	2.09	0.64
1:A:119:GLN:NE2	6:A:401:HOH:O	2.31	0.62
1:B:55:ARG:NH2	2:E:10:G:H4'	2.10	0.61
1:A:5:HIS:HE1	1:A:100:GLU:OE2	1.86	0.59
2:E:9:C:H6	2:E:9:C:H5"	1.69	0.57
1:B:52:ASN:ND2	1:B:54:LYS:HB3	2.20	0.57
1:B:52:ASN:OD1	1:B:53:GLU:N	2.38	0.56
1:A:154:ARG:NH1	6:A:404:HOH:O	2.40	0.54
1:B:154:ARG:NE	2:E:5:U:O4	2.42	0.52
1:A:91:THR:HG22	1:A:92:TRP:HD1	1.77	0.49
1:A:235:PHE:HB2	5:A:301:TLA:O1	2.11	0.49
1:A:199:VAL:HG11	2:D:-1:C:O2	2.14	0.48
2:D:12:G:C5'	2:D:12:G:H8	2.28	0.47
1:B:33:PRO:HB3	1:B:38:ARG:NH1	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:11:C:C2'	2:D:12:G:H5"	2.36	0.47
2:E:9:C:H5"	2:E:9:C:C6	2.50	0.46
1:A:61:GLU:O	1:A:62:GLN:HG2	2.17	0.45
1:A:85:ARG:HG2	1:A:241:SER:HA	1.99	0.45
1:B:52:ASN:O	1:B:56:VAL:HG23	2.19	0.42
1:B:38:ARG:NH2	1:B:43:GLY:O	2.52	0.42
1:A:17:LEU:O	1:A:93:VAL:HG11	2.19	0.42
1:A:29:HIS:HD2	6:A:524:HOH:O	2.03	0.41
1:B:85:ARG:HG2	1:B:241:SER:HA	2.02	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	249/266 (94%)	248 (100%)	1 (0%)	0	100	100
1	B	248/266 (93%)	246 (99%)	2 (1%)	0	100	100
All	All	497/532 (93%)	494 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	207/219 (94%)	205 (99%)	2 (1%)	76	63
1	B	206/219 (94%)	201 (98%)	5 (2%)	49	26
All	All	413/438 (94%)	406 (98%)	7 (2%)	60	42

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	50	MET
1	B	53	GLU
1	B	137	LYS
1	B	149	ASN
1	B	185	LEU
1	A	91	THR
1	A	149	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	251	HIS
1	A	5	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	D	20/26 (76%)	3 (15%)	1 (5%)
2	E	18/26 (69%)	0	0
All	All	38/52 (73%)	3 (7%)	1 (2%)

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	D	0	G
2	D	7	C
2	D	12	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	D	12	G

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

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	23G	D	22	2	19,29,30	5.47	14 (73%)	21,45,48	4.14	9 (42%)
2	23G	E	22	2	19,29,30	5.45	15 (78%)	21,45,48	3.86	10 (47%)

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
2	23G	D	22	2	-	0/3/35/36	0/4/4/4
2	23G	E	22	2	-	0/3/35/36	0/4/4/4

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	22	23G	C3'-C2'	-11.48	1.28	1.53
2	E	22	23G	C3'-C2'	-11.45	1.28	1.53
2	E	22	23G	C4-N3	9.54	1.50	1.35
2	D	22	23G	C4-N3	9.25	1.50	1.35
2	E	22	23G	C2-N2	7.78	1.49	1.33
2	D	22	23G	C2-N2	7.63	1.49	1.33
2	D	22	23G	C6-N1	7.41	1.45	1.33
2	D	22	23G	C6-C5	7.34	1.54	1.41
2	E	22	23G	C6-N1	7.08	1.45	1.33
2	E	22	23G	C6-C5	7.00	1.53	1.41
2	E	22	23G	O3'-C3'	6.38	1.59	1.45
2	D	22	23G	O3'-C3'	6.34	1.59	1.45
2	D	22	23G	O4'-C4'	-5.90	1.31	1.45
2	E	22	23G	O4'-C4'	-5.85	1.31	1.45
2	E	22	23G	O2'-C2'	5.63	1.57	1.45
2	D	22	23G	O2'-C2'	5.47	1.57	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	22	23G	C2-N1	4.69	1.43	1.35
2	D	22	23G	O4'-C1'	4.57	1.47	1.41
2	E	22	23G	O4'-C1'	4.45	1.47	1.41
2	E	22	23G	C2-N1	4.11	1.42	1.35
2	E	22	23G	C5-C4	-2.96	1.33	1.40
2	D	22	23G	C3'-C4'	2.94	1.60	1.52
2	E	22	23G	C3'-C4'	2.86	1.60	1.52
2	D	22	23G	C5-C4	-2.77	1.33	1.40
2	E	22	23G	O5'-C5'	-2.68	1.38	1.44
2	D	22	23G	O5'-C5'	-2.36	1.39	1.44
2	E	22	23G	PC-OC1	2.35	1.66	1.55
2	E	22	23G	C2-N3	2.24	1.45	1.34
2	D	22	23G	C2-N3	2.22	1.45	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	22	23G	N3-C2-N1	-9.99	113.89	127.22
2	E	22	23G	N3-C2-N1	-9.65	114.35	127.22
2	D	22	23G	C1'-N9-C4	7.69	140.15	126.64
2	D	22	23G	N2-C2-N1	7.26	128.55	117.25
2	E	22	23G	C1'-N9-C4	7.05	139.03	126.64
2	D	22	23G	C2-N3-C4	6.50	122.78	115.36
2	E	22	23G	C2-N3-C4	5.86	122.05	115.36
2	E	22	23G	N2-C2-N1	5.73	126.16	117.25
2	E	22	23G	C6-N1-C2	5.35	124.43	115.93
2	D	22	23G	C6-N1-C2	5.23	124.23	115.93
2	D	22	23G	C6-C5-C4	-4.81	116.21	120.80
2	D	22	23G	C4-C5-N7	-4.70	104.50	109.40
2	E	22	23G	C6-C5-C4	-4.32	116.67	120.80
2	E	22	23G	O3'-PC-OC2	-3.75	105.87	115.76
2	E	22	23G	C4-C5-N7	-3.56	105.69	109.40
2	E	22	23G	C5-C6-N1	-3.16	119.11	123.43
2	D	22	23G	C5-C6-N1	-2.99	119.35	123.43
2	E	22	23G	C5'-C4'-C3'	-2.15	107.27	114.40
2	D	22	23G	O5'-C5'-C4'	-2.08	101.92	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	TLA	E	101	-	3,9,9	0.72	0	6,12,12	1.30	1 (16%)
5	TLA	A	301	-	3,9,9	1.24	0	6,12,12	3.08	3 (50%)
5	TLA	A	302	-	3,9,9	0.91	0	6,12,12	1.26	1 (16%)
3	GOL	B	301	-	5,5,5	1.06	0	5,5,5	0.91	0
5	TLA	D	101	-	3,9,9	0.77	0	6,12,12	1.41	2 (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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	TLA	E	101	-	-	0/4/12/12	-
5	TLA	A	301	-	-	4/4/12/12	-
5	TLA	A	302	-	-	0/4/12/12	-
3	GOL	B	301	-	-	0/4/4/4	-
5	TLA	D	101	-	-	0/4/12/12	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	301	TLA	C1-C2-C3	-6.56	98.99	113.11
5	D	101	TLA	C4-C3-C2	-2.55	107.61	113.11
5	A	301	TLA	O3-C3-C4	2.44	116.97	111.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	E	101	TLA	C1-C2-C3	-2.40	107.94	113.11
5	A	301	TLA	C4-C3-C2	-2.38	107.98	113.11
5	A	302	TLA	O2-C2-C1	-2.23	105.73	111.10
5	D	101	TLA	C1-C2-C3	-2.10	108.58	113.11

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	301	TLA	C1-C2-C3-O3
5	A	301	TLA	C1-C2-C3-C4
5	A	301	TLA	O2-C2-C3-O3
5	A	301	TLA	O2-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	301	TLA	3	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	250/266 (93%)	0.23	12 (4%) 30 36	26, 36, 52, 75	0
1	B	250/266 (93%)	0.47	16 (6%) 19 25	24, 38, 57, 93	0
2	D	22/26 (84%)	-0.04	0 100 100	32, 42, 50, 59	0
2	E	23/26 (88%)	0.11	2 (8%) 10 13	33, 43, 48, 88	0
All	All	545/584 (93%)	0.33	30 (5%) 25 31	24, 37, 56, 93	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	ARG	9.7
1	B	55	ARG	5.2
1	B	56	VAL	5.2
1	B	54	LYS	5.1
1	B	53	GLU	4.7
1	A	57	ARG	4.5
1	B	50	MET	4.4
1	B	123	ARG	3.7
1	B	57	ARG	3.6
1	A	62	GLN	3.4
1	B	49	ALA	3.4
1	B	251	HIS	3.2
1	A	53	GLU	3.1
1	B	52	ASN	3.0
1	A	54	LYS	3.0
2	E	-3	C	2.8
1	A	50	MET	2.6
1	A	123	ARG	2.6
1	B	209	SER	2.5
1	A	75	ARG	2.3
1	A	154	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	30	ASP	2.3
1	B	14	GLY	2.2
1	A	251	HIS	2.2
1	B	64	GLY	2.2
1	B	210	THR	2.2
2	E	-1[A]	C	2.2
1	B	30	ASP	2.1
1	A	61	GLU	2.1
1	A	64	GLY	2.1

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
2	23G	E	22	26/27	0.85	0.13	30,30,30,30	0
2	23G	D	22	26/27	0.93	0.09	30,30,30,30	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. 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
5	TLA	A	302	10/10	0.64	0.25	55,83,98,110	0
5	TLA	E	101	10/10	0.85	0.12	49,56,64,71	0
5	TLA	D	101	10/10	0.90	0.12	51,58,60,62	0
3	GOL	B	301	6/6	0.94	0.17	46,51,55,66	0
5	TLA	A	301	10/10	0.94	0.30	44,53,82,91	0
4	MN	B	302	1/1	0.99	0.17	40,40,40,40	0

6.5 Other polymers (i)

There are no such residues in this entry.