



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 15, 2023 – 08:42 PM EDT

PDB ID : 8E2Q
Title : Crystal structure of TadAC-1.17 in a complex with ssDNA
Authors : Feliciano, P.R.; Lee, S.J.; Ciaramella, G.
Deposited on : 2022-08-15
Resolution : 2.34 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

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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (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

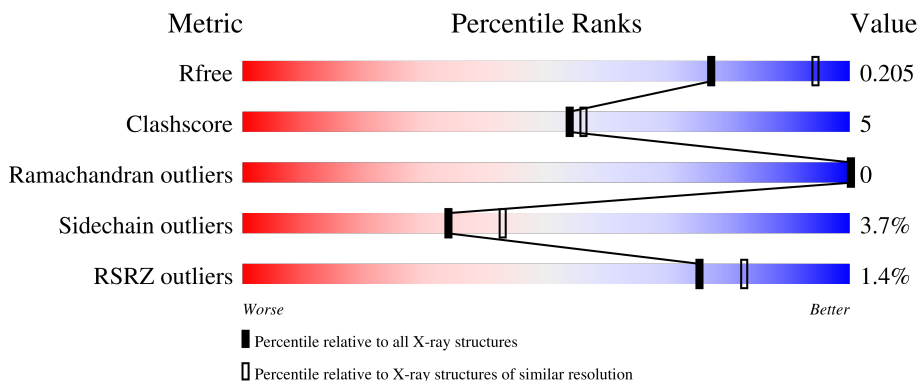
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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.



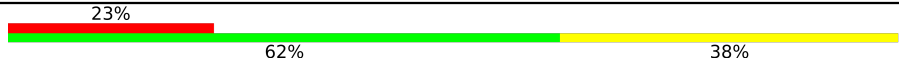
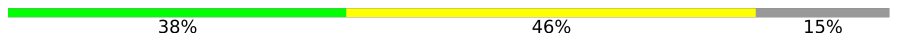

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	167	
1	B	167	
1	C	167	
1	D	167	
2	E	13	

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Mol	Chain	Length	Quality of chain
2	F	13	 <p>23% 62% 38%</p>
2	G	13	 <p>38% 46% 15%</p>
2	H	13	 <p>15% 38% 54% 8%</p>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5989 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 tRNA-specific adenosine deaminase 1.17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	154	1229	772	234	211	12	0	3	0
1	B	160	1265	793	240	219	13	0	1	0
1	C	153	1210	763	227	207	13	0	1	0
1	D	154	1162	735	211	204	12	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	ALA	THR	conflict	UNP W8T8U5
A	23	ARG	TRP	conflict	UNP W8T8U5
A	36	LEU	HIS	conflict	UNP W8T8U5
A	48	GLY	PRO	conflict	UNP W8T8U5
A	51	LEU	ARG	conflict	UNP W8T8U5
A	76	TYR	ILE	conflict	UNP W8T8U5
A	82	THR	VAL	conflict	UNP W8T8U5
A	84	PHE	LEU	conflict	UNP W8T8U5
A	106	VAL	ALA	conflict	UNP W8T8U5
A	108	ASN	ASP	conflict	UNP W8T8U5
A	142	GLU	ALA	conflict	UNP W8T8U5
A	146	CYS	SER	conflict	UNP W8T8U5
A	147	ARG	ASP	conflict	UNP W8T8U5
A	152	PRO	ARG	conflict	UNP W8T8U5
A	154	ARG	GLN	conflict	UNP W8T8U5
A	155	VAL	GLU	conflict	UNP W8T8U5
A	156	PHE	ILE	conflict	UNP W8T8U5
A	157	ASN	LYS	conflict	UNP W8T8U5
B	17	ALA	THR	conflict	UNP W8T8U5
B	23	ARG	TRP	conflict	UNP W8T8U5
B	36	LEU	HIS	conflict	UNP W8T8U5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	48	GLY	PRO	conflict	UNP W8T8U5
B	51	LEU	ARG	conflict	UNP W8T8U5
B	76	TYR	ILE	conflict	UNP W8T8U5
B	82	THR	VAL	conflict	UNP W8T8U5
B	84	PHE	LEU	conflict	UNP W8T8U5
B	106	VAL	ALA	conflict	UNP W8T8U5
B	108	ASN	ASP	conflict	UNP W8T8U5
B	142	GLU	ALA	conflict	UNP W8T8U5
B	146	CYS	SER	conflict	UNP W8T8U5
B	147	ARG	ASP	conflict	UNP W8T8U5
B	152	PRO	ARG	conflict	UNP W8T8U5
B	154	ARG	GLN	conflict	UNP W8T8U5
B	155	VAL	GLU	conflict	UNP W8T8U5
B	156	PHE	ILE	conflict	UNP W8T8U5
B	157	ASN	LYS	conflict	UNP W8T8U5
C	17	ALA	THR	conflict	UNP W8T8U5
C	23	ARG	TRP	conflict	UNP W8T8U5
C	36	LEU	HIS	conflict	UNP W8T8U5
C	48	GLY	PRO	conflict	UNP W8T8U5
C	51	LEU	ARG	conflict	UNP W8T8U5
C	76	TYR	ILE	conflict	UNP W8T8U5
C	82	THR	VAL	conflict	UNP W8T8U5
C	84	PHE	LEU	conflict	UNP W8T8U5
C	106	VAL	ALA	conflict	UNP W8T8U5
C	108	ASN	ASP	conflict	UNP W8T8U5
C	142	GLU	ALA	conflict	UNP W8T8U5
C	146	CYS	SER	conflict	UNP W8T8U5
C	147	ARG	ASP	conflict	UNP W8T8U5
C	152	PRO	ARG	conflict	UNP W8T8U5
C	154	ARG	GLN	conflict	UNP W8T8U5
C	155	VAL	GLU	conflict	UNP W8T8U5
C	156	PHE	ILE	conflict	UNP W8T8U5
C	157	ASN	LYS	conflict	UNP W8T8U5
D	17	ALA	THR	conflict	UNP W8T8U5
D	23	ARG	TRP	conflict	UNP W8T8U5
D	36	LEU	HIS	conflict	UNP W8T8U5
D	48	GLY	PRO	conflict	UNP W8T8U5
D	51	LEU	ARG	conflict	UNP W8T8U5
D	76	TYR	ILE	conflict	UNP W8T8U5
D	82	THR	VAL	conflict	UNP W8T8U5
D	84	PHE	LEU	conflict	UNP W8T8U5
D	106	VAL	ALA	conflict	UNP W8T8U5

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Chain	Residue	Modelled	Actual	Comment	Reference
D	108	ASN	ASP	conflict	UNP W8T8U5
D	142	GLU	ALA	conflict	UNP W8T8U5
D	146	CYS	SER	conflict	UNP W8T8U5
D	147	ARG	ASP	conflict	UNP W8T8U5
D	152	PRO	ARG	conflict	UNP W8T8U5
D	154	ARG	GLN	conflict	UNP W8T8U5
D	155	VAL	GLU	conflict	UNP W8T8U5
D	156	PHE	ILE	conflict	UNP W8T8U5
D	157	ASN	LYS	conflict	UNP W8T8U5

- Molecule 2 is a DNA chain called DNA (5'-D(P*GP*CP*GP*GP*CP*TP*(D8A)P*CP*GP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	11	Total 229	C 106	N 46	O 66	P 11	0	0	0
2	F	13	Total 268	C 125	N 51	O 79	P 13	0	0	0
2	G	11	Total 227	C 106	N 43	O 67	P 11	0	0	0
2	H	12	Total 246	C 115	N 46	O 73	P 12	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Zn 1	0	0
3	B	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	D	1	Total 1	Zn 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is water.

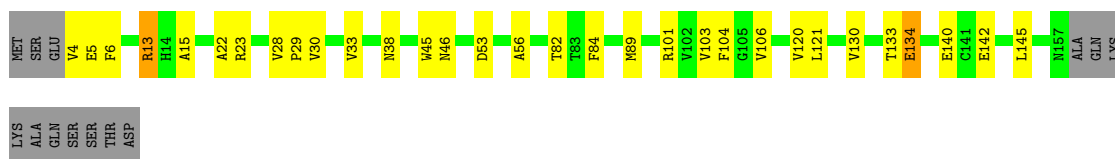
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	30	Total O 30 30	0	0
5	B	27	Total O 27 27	0	0
5	C	24	Total O 24 24	0	0
5	D	26	Total O 26 26	0	0
5	E	3	Total O 3 3	0	0
5	F	7	Total O 7 7	0	0
5	G	6	Total O 6 6	0	0
5	H	8	Total O 8 8	0	0

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: tRNA-specific adenosine deaminase 1.17

Chain A: 




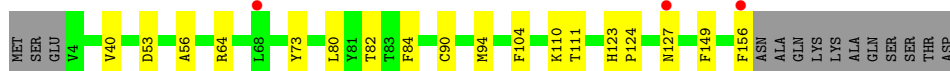
- Molecule 1: tRNA-specific adenosine deaminase 1.17

Chain B: 




- Molecule 1: tRNA-specific adenosine deaminase 1.17

Chain C: 



- Molecule 1: tRNA-specific adenosine deaminase 1.17

Chain D: 

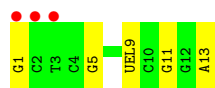


- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*CP*TP*(D8A)P*CP*GP*GP*A)-3')

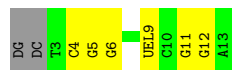
Chain E: 



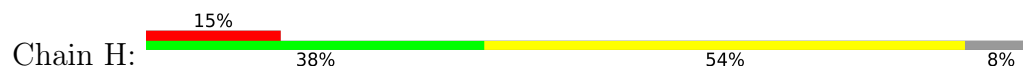
- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*CP*TP*(D8A)P*CP*GP*GP*A)-3')



- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*CP*TP*(D8A)P*CP*GP*GP*A)-3')



- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*CP*TP*(D8A)P*CP*GP*GP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.93Å 85.93Å 224.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	62.03 – 2.34 62.03 – 2.34	Depositor EDS
% Data completeness (in resolution range)	100.0 (62.03-2.34) 100.0 (62.03-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	215.91 (at 2.34Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.169 , 0.205 0.170 , 0.205	Depositor DCC
R_{free} test set	2084 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	46.4	Xtrriage
Anisotropy	0.341	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 30.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	0.230 for -h,-k,l	Xtrriage
Reported twinning fraction	0.280 for -h,-k,l	Depositor
Outliers	0 of 41494 reflections	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5989	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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: UEL, ZN, GOL

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	0.27	0/1255	0.51	0/1695
1	B	0.28	0/1291	0.53	0/1741
1	C	0.26	0/1236	0.52	0/1668
1	D	0.27	0/1187	0.50	0/1608
2	E	0.42	0/231	0.77	0/350
2	F	0.49	0/275	0.82	0/420
2	G	0.53	0/229	0.86	0/349
2	H	0.57	0/250	0.91	0/381
All	All	0.32	0/5954	0.59	0/8212

There are no bond length outliers.

There are no bond angle outliers.

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	1229	0	1207	20	0
1	B	1265	0	1251	15	0
1	C	1210	0	1199	10	0
1	D	1162	0	1113	11	0
2	E	229	0	113	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	268	0	135	4	0
2	G	227	0	113	3	0
2	H	246	0	124	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	6	0	8	0	0
4	C	6	0	8	0	0
4	D	6	0	8	1	0
5	A	30	0	0	3	0
5	B	27	0	0	1	0
5	C	24	0	0	0	0
5	D	26	0	0	1	0
5	E	3	0	0	0	0
5	F	7	0	0	0	0
5	G	6	0	0	0	0
5	H	8	0	0	0	0
All	All	5989	0	5279	59	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 5.

The worst 5 of 59 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:C:110:LYS:HD3	1:C:111:THR:HG23	1.73	0.71
1:A:22:ALA:HB2	1:A:45:TRP:HB2	1.78	0.66
2:G:4:DC:H2''	2:G:5:DG:H5'	1.84	0.59
1:A:101[A]:ARG:NH1	5:A:302:HOH:O	2.37	0.58
1:A:82:THR:O	1:A:104:PHE:HA	2.04	0.56

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	155/167 (93%)	148 (96%)	7 (4%)	0	100	100
1	B	159/167 (95%)	154 (97%)	5 (3%)	0	100	100
1	C	152/167 (91%)	146 (96%)	6 (4%)	0	100	100
1	D	152/167 (91%)	145 (95%)	7 (5%)	0	100	100
All	All	618/668 (92%)	593 (96%)	25 (4%)	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	124/135 (92%)	118 (95%)	6 (5%)	25	32
1	B	129/135 (96%)	126 (98%)	3 (2%)	50	61
1	C	123/135 (91%)	118 (96%)	5 (4%)	30	38
1	D	112/135 (83%)	107 (96%)	5 (4%)	27	34
All	All	488/540 (90%)	469 (96%)	19 (4%)	34	41

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	55	THR
1	D	84	PHE
1	D	97	SER
1	D	74	ARG
1	B	84	PHE

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

Mol	Chain	Res	Type
1	B	122	HIS
1	C	96	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

4 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	UEL	F	9	3,2	16,23,24	4.91	11 (68%)	14,33,36	1.77	4 (28%)
2	UEL	G	9	3,2	16,23,24	4.89	10 (62%)	14,33,36	1.48	3 (21%)
2	UEL	E	9	3,2	16,23,24	4.91	11 (68%)	14,33,36	1.68	3 (21%)
2	UEL	H	9	3,2	16,23,24	4.92	10 (62%)	14,33,36	1.73	3 (21%)

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	UEL	F	9	3,2	-	2/3/31/32	0/3/3/3
2	UEL	G	9	3,2	-	2/3/31/32	0/3/3/3
2	UEL	E	9	3,2	-	2/3/31/32	0/3/3/3
2	UEL	H	9	3,2	-	2/3/31/32	0/3/3/3

The worst 5 of 42 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	9	UEL	N7-N8	-11.13	1.14	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	9	UEL	N7-N8	-11.12	1.14	1.34
2	E	9	UEL	N7-N8	-11.09	1.14	1.34
2	G	9	UEL	N7-N8	-10.96	1.14	1.34
2	G	9	UEL	N8-N9	-9.63	1.17	1.34

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	9	UEL	C2'-C1'-N9	-4.55	108.31	113.67
2	E	9	UEL	C2'-C1'-N9	-4.27	108.63	113.67
2	F	9	UEL	C2'-C1'-N9	-3.59	109.44	113.67
2	G	9	UEL	C2'-C1'-N9	-3.44	109.62	113.67
2	F	9	UEL	C4'-O4'-C1'	-3.12	101.92	109.45

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	9	UEL	O4'-C4'-C5'-O5'
2	G	9	UEL	O4'-C4'-C5'-O5'
2	H	9	UEL	O4'-C4'-C5'-O5'
2	E	9	UEL	C3'-C4'-C5'-O5'
2	F	9	UEL	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 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	GOL	D	201	-	5,5,5	0.96	0	5,5,5	1.00	0
4	GOL	B	201	-	5,5,5	0.87	0	5,5,5	1.08	0
4	GOL	C	201	-	5,5,5	0.93	0	5,5,5	1.01	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	GOL	D	201	-	-	0/4/4/4	-
4	GOL	B	201	-	-	2/4/4/4	-
4	GOL	C	201	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	201	GOL	C1-C2-C3-O3
4	B	201	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	201	GOL	1	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	154/167 (92%)	0.01	0 100 100	29, 37, 53, 76	0
1	B	160/167 (95%)	0.06	0 100 100	28, 36, 53, 76	0
1	C	153/167 (91%)	0.10	3 (1%) 65 74	30, 40, 58, 75	0
1	D	154/167 (92%)	0.22	1 (0%) 89 93	33, 45, 65, 81	0
2	E	10/13 (76%)	-0.10	0 100 100	41, 45, 55, 124	0
2	F	12/13 (92%)	0.45	3 (25%) 0 1	37, 45, 107, 123	0
2	G	10/13 (76%)	-0.28	0 100 100	36, 41, 45, 67	0
2	H	11/13 (84%)	0.15	2 (18%) 1 2	37, 40, 65, 77	0
All	All	664/720 (92%)	0.09	9 (1%) 75 82	28, 40, 63, 124	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	3	DT	4.4
1	C	156	PHE	3.7
1	D	127	ASN	3.5
2	F	2	DC	3.4
2	F	3	DT	3.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	UEL	E	9	21/22	0.95	0.16	32,38,44,49	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	UEL	F	9	21/22	0.95	0.13	30,33,40,41	0
2	UEL	G	9	21/22	0.95	0.13	35,41,50,56	0
2	UEL	H	9	21/22	0.96	0.13	38,42,45,53	0

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, 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(\AA^2)	Q<0.9
4	GOL	B	201	6/6	0.86	0.23	35,39,42,47	0
4	GOL	C	201	6/6	0.87	0.18	43,44,48,52	0
4	GOL	D	201	6/6	0.87	0.20	54,56,58,61	0
3	ZN	C	202	1/1	0.98	0.14	42,42,42,42	0
3	ZN	B	202	1/1	0.99	0.17	38,38,38,38	0
3	ZN	A	201	1/1	0.99	0.16	38,38,38,38	0
3	ZN	D	202	1/1	0.99	0.14	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.