



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 07:03 AM GMT

PDB ID : 3N4Q
Title : Human cytomegalovirus terminase nuclease domain, Mn soaked
Authors : Nadal, M.; Mas, P.J.; Blanco, A.G.; Arnan, C.; Sola, M.; Hart, D.J.; Coll, M.
Deposited on : 2010-05-22
Resolution : 3.20 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

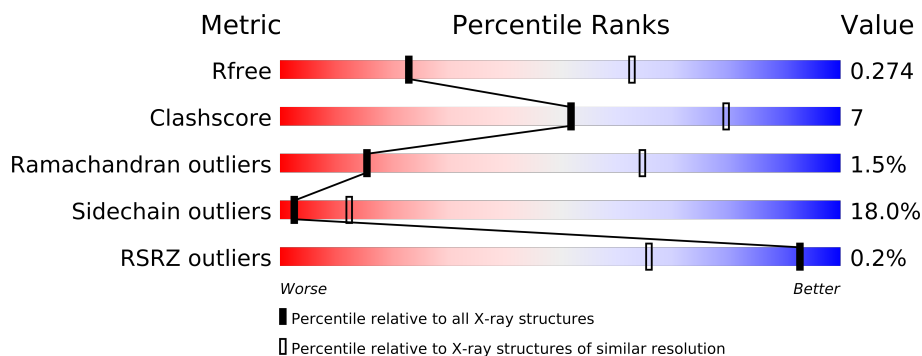
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.20 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}	66092	1824 (3.30-3.10)
Clashscore	79885	1078 (3.26-3.14)
Ramachandran outliers	78287	1059 (3.26-3.14)
Sidechain outliers	78261	1058 (3.26-3.14)
RSRZ outliers	66119	1825 (3.30-3.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	279	
1	B	279	
1	C	279	
1	D	279	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7056 atoms, of which 0 are hydrogen and 0 are deuterium.

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 TERMINASE SUBUNIT UL89 PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1736	1113	292	325	6			
1	B	221	Total	C	N	O	S	0	0	0
			1775	1137	301	331	6			
1	C	219	Total	C	N	O	S	0	0	0
			1755	1126	293	330	6			
1	D	214	Total	C	N	O	S	0	0	0
			1722	1104	290	322	6			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	396	MET	-	EXPRESSION TAG	UNP P16732
A	397	GLY	-	EXPRESSION TAG	UNP P16732
A	398	HIS	-	EXPRESSION TAG	UNP P16732
A	399	HIS	-	EXPRESSION TAG	UNP P16732
A	400	HIS	-	EXPRESSION TAG	UNP P16732
A	401	HIS	-	EXPRESSION TAG	UNP P16732
A	402	HIS	-	EXPRESSION TAG	UNP P16732
A	403	HIS	-	EXPRESSION TAG	UNP P16732
A	404	ASP	-	EXPRESSION TAG	UNP P16732
A	405	TYR	-	EXPRESSION TAG	UNP P16732
A	406	ASP	-	EXPRESSION TAG	UNP P16732
A	407	ILE	-	EXPRESSION TAG	UNP P16732
A	408	PRO	-	EXPRESSION TAG	UNP P16732
A	409	THR	-	EXPRESSION TAG	UNP P16732
A	410	THR	-	EXPRESSION TAG	UNP P16732
A	411	GLU	-	EXPRESSION TAG	UNP P16732
A	412	ASN	-	EXPRESSION TAG	UNP P16732
A	413	LEU	-	EXPRESSION TAG	UNP P16732
A	414	TYR	-	EXPRESSION TAG	UNP P16732
A	415	PHE	-	EXPRESSION TAG	UNP P16732
A	416	GLN	-	EXPRESSION TAG	UNP P16732

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Chain	Residue	Modelled	Actual	Comment	Reference
A	417	GLY	-	EXPRESSION TAG	UNP P16732
B	396	MET	-	EXPRESSION TAG	UNP P16732
B	397	GLY	-	EXPRESSION TAG	UNP P16732
B	398	HIS	-	EXPRESSION TAG	UNP P16732
B	399	HIS	-	EXPRESSION TAG	UNP P16732
B	400	HIS	-	EXPRESSION TAG	UNP P16732
B	401	HIS	-	EXPRESSION TAG	UNP P16732
B	402	HIS	-	EXPRESSION TAG	UNP P16732
B	403	HIS	-	EXPRESSION TAG	UNP P16732
B	404	ASP	-	EXPRESSION TAG	UNP P16732
B	405	TYR	-	EXPRESSION TAG	UNP P16732
B	406	ASP	-	EXPRESSION TAG	UNP P16732
B	407	ILE	-	EXPRESSION TAG	UNP P16732
B	408	PRO	-	EXPRESSION TAG	UNP P16732
B	409	THR	-	EXPRESSION TAG	UNP P16732
B	410	THR	-	EXPRESSION TAG	UNP P16732
B	411	GLU	-	EXPRESSION TAG	UNP P16732
B	412	ASN	-	EXPRESSION TAG	UNP P16732
B	413	LEU	-	EXPRESSION TAG	UNP P16732
B	414	TYR	-	EXPRESSION TAG	UNP P16732
B	415	PHE	-	EXPRESSION TAG	UNP P16732
B	416	GLN	-	EXPRESSION TAG	UNP P16732
B	417	GLY	-	EXPRESSION TAG	UNP P16732
C	396	MET	-	EXPRESSION TAG	UNP P16732
C	397	GLY	-	EXPRESSION TAG	UNP P16732
C	398	HIS	-	EXPRESSION TAG	UNP P16732
C	399	HIS	-	EXPRESSION TAG	UNP P16732
C	400	HIS	-	EXPRESSION TAG	UNP P16732
C	401	HIS	-	EXPRESSION TAG	UNP P16732
C	402	HIS	-	EXPRESSION TAG	UNP P16732
C	403	HIS	-	EXPRESSION TAG	UNP P16732
C	404	ASP	-	EXPRESSION TAG	UNP P16732
C	405	TYR	-	EXPRESSION TAG	UNP P16732
C	406	ASP	-	EXPRESSION TAG	UNP P16732
C	407	ILE	-	EXPRESSION TAG	UNP P16732
C	408	PRO	-	EXPRESSION TAG	UNP P16732
C	409	THR	-	EXPRESSION TAG	UNP P16732
C	410	THR	-	EXPRESSION TAG	UNP P16732
C	411	GLU	-	EXPRESSION TAG	UNP P16732
C	412	ASN	-	EXPRESSION TAG	UNP P16732
C	413	LEU	-	EXPRESSION TAG	UNP P16732
C	414	TYR	-	EXPRESSION TAG	UNP P16732

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Chain	Residue	Modelled	Actual	Comment	Reference
C	415	PHE	-	EXPRESSION TAG	UNP P16732
C	416	GLN	-	EXPRESSION TAG	UNP P16732
C	417	GLY	-	EXPRESSION TAG	UNP P16732
D	396	MET	-	EXPRESSION TAG	UNP P16732
D	397	GLY	-	EXPRESSION TAG	UNP P16732
D	398	HIS	-	EXPRESSION TAG	UNP P16732
D	399	HIS	-	EXPRESSION TAG	UNP P16732
D	400	HIS	-	EXPRESSION TAG	UNP P16732
D	401	HIS	-	EXPRESSION TAG	UNP P16732
D	402	HIS	-	EXPRESSION TAG	UNP P16732
D	403	HIS	-	EXPRESSION TAG	UNP P16732
D	404	ASP	-	EXPRESSION TAG	UNP P16732
D	405	TYR	-	EXPRESSION TAG	UNP P16732
D	406	ASP	-	EXPRESSION TAG	UNP P16732
D	407	ILE	-	EXPRESSION TAG	UNP P16732
D	408	PRO	-	EXPRESSION TAG	UNP P16732
D	409	THR	-	EXPRESSION TAG	UNP P16732
D	410	THR	-	EXPRESSION TAG	UNP P16732
D	411	GLU	-	EXPRESSION TAG	UNP P16732
D	412	ASN	-	EXPRESSION TAG	UNP P16732
D	413	LEU	-	EXPRESSION TAG	UNP P16732
D	414	TYR	-	EXPRESSION TAG	UNP P16732
D	415	PHE	-	EXPRESSION TAG	UNP P16732
D	416	GLN	-	EXPRESSION TAG	UNP P16732
D	417	GLY	-	EXPRESSION TAG	UNP P16732

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Mn 2 2	0	0
2	A	2	Total Mn 2 2	0	0
2	D	2	Total Mn 2 2	0	0
2	C	2	Total Mn 2 2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	11	Total 11	O 11	0	0
4	B	17	Total 17	O 17	0	0
4	C	15	Total 15	O 15	0	0
4	D	13	Total 13	O 13	0	0

E622	Q625	Q626	L627	H628	R629	VAL	THR	LEU	THR	ALA	GLU	GLY	THR	THR	ALA	ARG	TYR	SER	ALA	ALA	LYS	ARG	GLN	ASN	ARG	ILE	SER	D650	L662	S674	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488	L492	F495	F496	LEU	ARG	ASP	LEU	LEU	SER	E502	S503	S504	A507	S518	S521	D527	E528	L529	N538	Q539	C547	L548	S552	F563	Y564	H565	N570	Q574	M579	G580	R581	I599	E604	S607	I610	S613	H614	L620	L621	H485	L488</
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.70Å 88.10Å 190.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.05 – 3.20 44.05 – 3.19	Depositor EDS
% Data completeness (in resolution range)	89.6 (44.05-3.20) 89.6 (44.05-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.90 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.207 , 0.283 0.201 , 0.274	Depositor DCC
R_{free} test set	1109 reflections (5.40%)	DCC
Wilson B-factor (Å ²)	75.1	Xtriage
Anisotropy	0.820	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 59.6	EDS
Estimated twinning fraction	0.046 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 21659 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7056	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.19% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.61	0/1770	0.75	1/2399 (0.0%)
1	B	0.58	1/1809 (0.1%)	0.74	2/2452 (0.1%)
1	C	0.58	0/1789	0.74	3/2427 (0.1%)
1	D	0.55	0/1756	0.69	0/2380
All	All	0.58	1/7124 (0.0%)	0.73	6/9658 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	622	GLU	CG-CD	5.10	1.59	1.51

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	488	LEU	CA-CB-CG	7.49	132.52	115.30
1	C	458	LEU	CA-CB-CG	6.16	129.46	115.30
1	B	605	LEU	CA-CB-CG	5.69	128.38	115.30
1	C	569	GLN	N-CA-C	-5.46	96.27	111.00
1	B	578	LEU	CA-CB-CG	5.35	127.60	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1736	0	0	9	0
1	B	1775	0	0	17	0
1	C	1755	0	0	12	0
1	D	1722	0	0	12	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	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	A	11	0	0	0	0
4	B	17	0	0	2	0
4	C	15	0	0	1	0
4	D	13	0	0	1	0
All	All	7056	0	0	50	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

The worst 5 of 50 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:538:ASN:OD1	1:A:541:ALA:N	2.15	0.79
1:B:664:ASP:CB	4:B:18:HOH:O	2.30	0.78
1:C:583:LYS:NZ	4:C:37:HOH:O	2.27	0.67
1:B:568:ASP:CG	1:B:569:GLN:N	2.49	0.65
1:A:432:ASP:OD1	1:A:432:ASP:N	2.32	0.62

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	208/279 (75%)	188 (90%)	15 (7%)	5 (2%)	9	51
1	B	213/279 (76%)	182 (85%)	29 (14%)	2 (1%)	25	76
1	C	213/279 (76%)	179 (84%)	30 (14%)	4 (2%)	12	59
1	D	206/279 (74%)	178 (86%)	26 (13%)	2 (1%)	22	74
All	All	840/1116 (75%)	727 (86%)	100 (12%)	13 (2%)	15	64

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	612	LEU
1	C	569	GLN
1	D	504	SER
1	A	537	THR
1	A	569	GLN

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	188/242 (78%)	154 (82%)	34 (18%)	2	12
1	B	192/242 (79%)	160 (83%)	32 (17%)	3	14
1	C	190/242 (78%)	154 (81%)	36 (19%)	2	11
1	D	186/242 (77%)	152 (82%)	34 (18%)	2	12
All	All	756/968 (78%)	620 (82%)	136 (18%)	2	12

5 of 136 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	663	CYS
1	C	529	LEU
1	D	579	MET
1	B	665	ASP
1	C	460	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 12 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

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	216/279 (77%)	-0.27	1 (0%) 88 46	48, 77, 119, 164	6 (2%)
1	B	221/279 (79%)	-0.33	0 100 100	53, 76, 110, 154	2 (0%)
1	C	219/279 (78%)	-0.26	1 (0%) 88 46	54, 82, 128, 180	4 (1%)
1	D	214/279 (76%)	-0.21	0 100 100	57, 86, 122, 148	12 (5%)
All	All	870/1116 (77%)	-0.27	2 (0%) 93 66	48, 80, 123, 180	24 (2%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	497	LEU	2.2
1	C	571	HIS	2.1

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MN	D	8	1/1	0.19	-0.31	64,64,64,64	0
3	MG	C	4	1/1	0.15	-1.09	65,65,65,65	0
2	MN	A	1	1/1	0.13	-1.17	58,58,58,58	0
2	MN	A	2	1/1	0.14	-1.31	55,55,55,55	0
2	MN	C	6	1/1	0.06	-1.73	58,58,58,58	0
3	MG	B	2	1/1	0.14	-1.84	40,40,40,40	0
2	MN	D	7	1/1	0.10	-1.94	60,60,60,60	0
2	MN	C	5	1/1	0.06	-2.18	58,58,58,58	0
3	MG	A	675	1/1	0.09	-2.21	29,29,29,29	0
3	MG	D	3	1/1	0.05	-3.26	38,38,38,38	0
2	MN	B	3	1/1	0.05	-3.61	55,55,55,55	0
2	MN	B	4	1/1	0.06	-6.31	52,52,52,52	0

6.5 Other polymers ⓘ

There are no such residues in this entry.