



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:30 PM GMT

PDB ID : 3ZNS
Title : HDAC7 bound with TFMO inhibitor tmp942
Authors : Lobera, M.; Madauss, K.P.; Pohlhaus, D.T.; Trump, R.P.; Nolan, M.A.
Deposited on : 2013-02-15
Resolution : 2.45 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

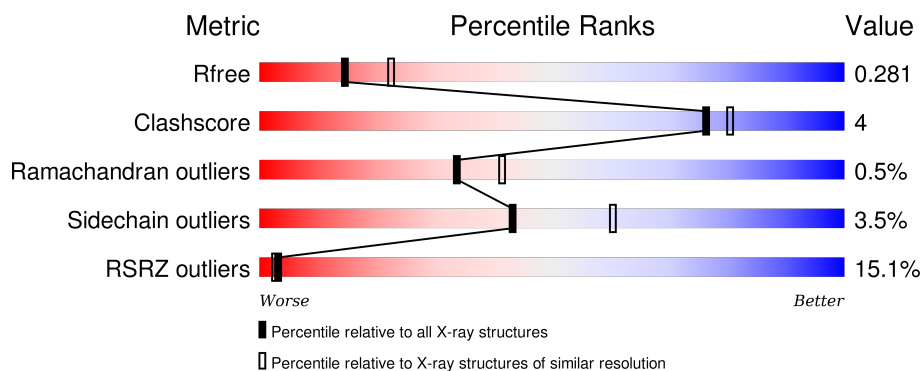
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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}	91344	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	423	<div> <div>10%</div> <div>78%</div> <div>8%</div> <div>13%</div> </div>
1	B	423	<div> <div>14%</div> <div>79%</div> <div>10%</div> <div>10%</div> </div>
1	C	423	<div> <div>15%</div> <div>76%</div> <div>9%</div> <div>15%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8601 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 HISTONE DEACETYLASE 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2774	1740	499	516	19			
1	B	380	Total	C	N	O	S	0	0	0
			2867	1798	514	536	19			
1	C	359	Total	C	N	O	S	0	0	0
			2704	1691	490	505	18			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	481	GLY	-	EXPRESSION TAG	UNP Q8WUI4
B	481	GLY	-	EXPRESSION TAG	UNP Q8WUI4
C	481	GLY	-	EXPRESSION TAG	UNP Q8WUI4

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

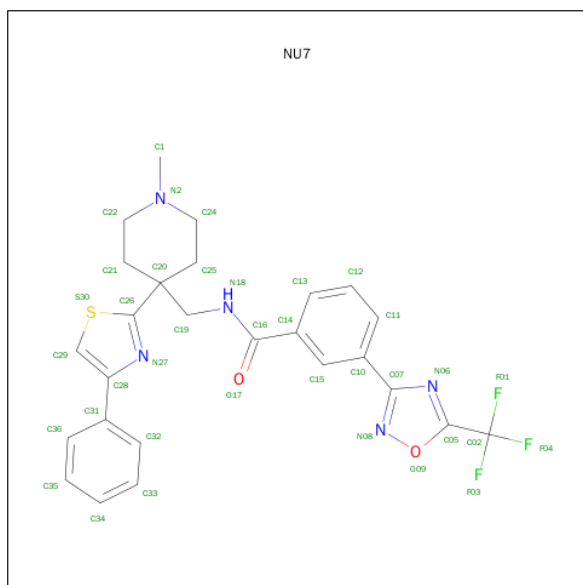
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	K	0	0
			2	2		
3	A	2	Total	K	0	0
			2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	2	Total K 2 2	0	0

- Molecule 4 is N-{[1-METHYL-4-(4-PHENYL-1,3-THIAZOL-2-YL)PIPERIDIN-4-YL]METHYL}-3-[5-(TRIFLUOROMETHYL)-1,2,4-OXADIAZOL-3-YL]BENZAMIDE (three-letter code: NU7) (formula: C₂₆H₂₄F₃N₅O₂S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C F N O S 37 26 3 5 2 1	0	0
4	B	1	Total C F N O S 37 26 3 5 2 1	0	0
4	C	1	Total C F N O S 37 26 3 5 2 1	0	0

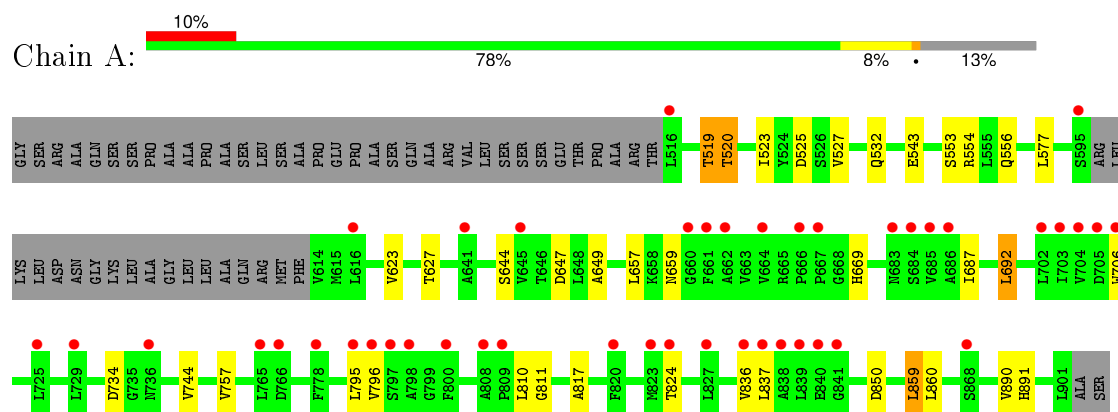
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	46	Total O 46 46	0	0
5	B	69	Total O 69 69	0	0
5	C	18	Total O 18 18	0	0

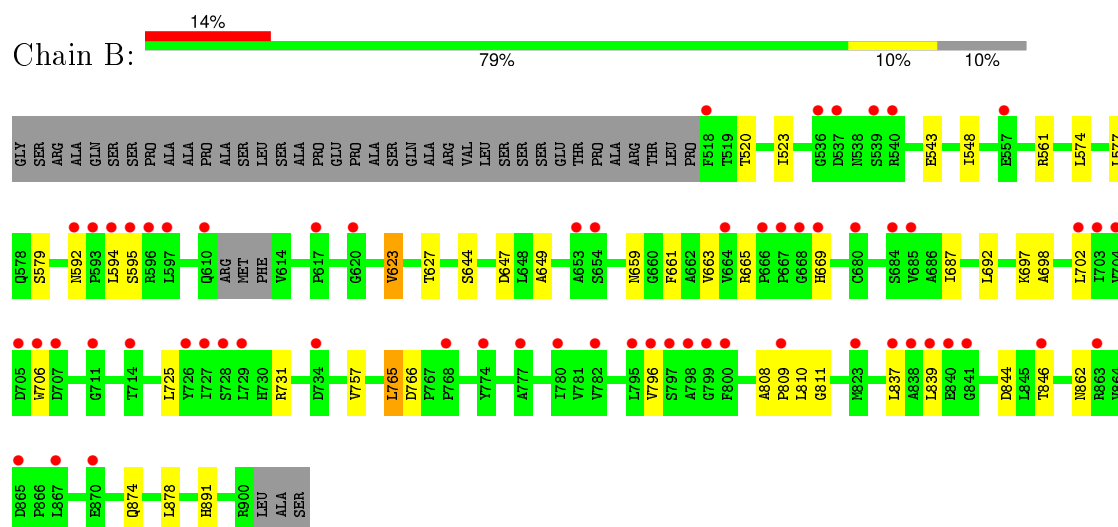
3 Residue-property plots

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

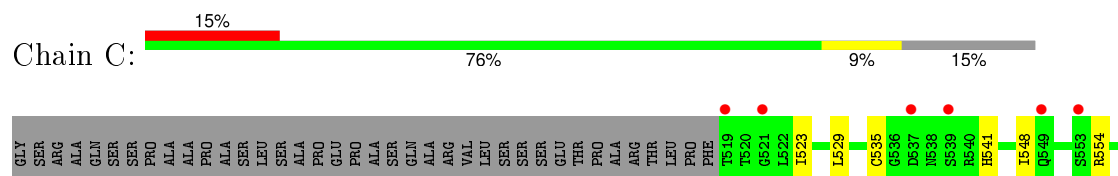
• Molecule 1: HISTONE DEACETYLASE 7

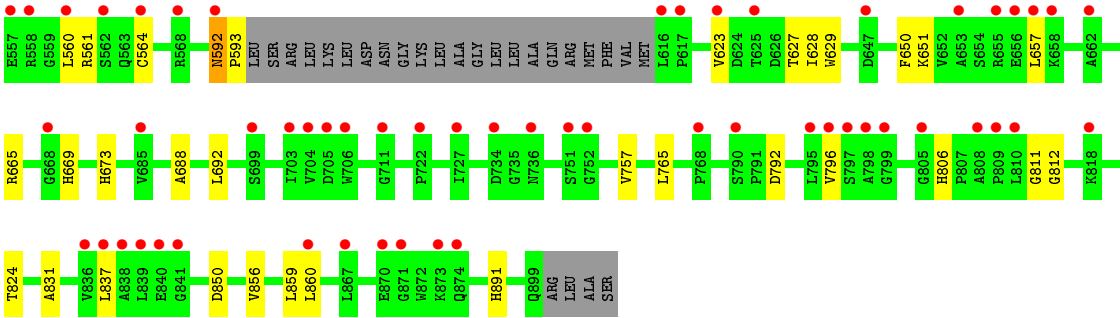


• Molecule 1: HISTONE DEACETYLASE 7



• Molecule 1: HISTONE DEACETYLASE 7





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	81.57Å 81.57Å 150.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	70.71 – 2.45 33.14 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.8 (70.71-2.45) 99.8 (33.14-2.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.88 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0053	Depositor
R, R_{free}	0.186 , 0.225 0.245 , 0.281	Depositor DCC
R_{free} test set	2055 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 30.0	EDS
Estimated twinning fraction	0.011 for -h,-k,l 0.045 for h,-h-k,-l 0.029 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 40864 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8601	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NU7, ZN, K

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	0.32	0/2842	0.49	0/3860
1	B	0.33	0/2934	0.48	0/3981
1	C	0.31	0/2770	0.47	0/3764
All	All	0.32	0/8546	0.48	0/11605

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

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	2774	0	2651	19	0
1	B	2867	0	2751	20	0
1	C	2704	0	2571	25	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
4	A	37	0	24	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	37	0	24	2	0
4	C	37	0	24	0	0
5	A	46	0	0	0	0
5	B	69	0	0	0	0
5	C	18	0	0	0	0
All	All	8601	0	8045	63	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 4.

All (63) 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:535:CYS:HG	2:C:102:ZN:ZN	0.69	0.94
1:C:623:VAL:HG22	1:C:627:THR:HB	1.66	0.77
1:B:623:VAL:HG22	1:B:627:THR:HB	1.67	0.76
1:B:523:ILE:HG21	1:B:644:SER:HB3	1.69	0.74
1:C:541:HIS:CE1	1:C:628:ILE:HD11	2.25	0.71
1:B:844:ASP:OD2	1:B:846:THR:HG22	1.92	0.68
1:A:623:VAL:HG22	1:A:627:THR:HB	1.76	0.65
1:B:808:ALA:HB3	1:B:809:PRO:HD3	1.80	0.63
1:C:796:VAL:HB	1:C:837:LEU:HD22	1.81	0.62
1:A:554:ARG:NH2	1:A:850:ASP:OD1	2.33	0.61
1:A:520:THR:HG23	1:A:860:LEU:HD23	1.82	0.61
1:C:651:LYS:HB3	1:C:657:LEU:HD13	1.84	0.60
1:C:592:ASN:CB	1:C:593:PRO:HA	2.33	0.59
1:A:744:VAL:HG12	1:A:890:VAL:HG21	1.85	0.58
1:A:649:ALA:HB1	1:A:692:LEU:HD21	1.86	0.57
1:C:688:ALA:O	1:C:692:LEU:HD13	2.05	0.57
1:C:548:ILE:HD13	1:C:665:ARG:HG2	1.88	0.55
1:B:796:VAL:HB	1:B:837:LEU:HD22	1.89	0.54
1:C:650:PHE:CE1	1:C:692:LEU:HD12	2.43	0.53
1:C:592:ASN:HB2	1:C:593:PRO:HA	1.91	0.53
1:C:564:CYS:SG	1:C:860:LEU:HD21	2.49	0.52
1:B:520:THR:HG22	1:B:659:ASN:OD1	2.09	0.52
1:B:810:LEU:HD22	4:B:1900:NU7:H29	1.92	0.52
1:A:519:THR:HG22	1:A:520:THR:H	1.74	0.52
1:B:661:PHE:CE2	1:B:663:VAL:HG22	2.46	0.50
1:B:649:ALA:HB1	1:B:692:LEU:HD21	1.92	0.49
1:C:806:HIS:O	1:C:812:GLY:N	2.38	0.49
1:C:796:VAL:CB	1:C:837:LEU:HD22	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:592:ASN:CB	1:C:593:PRO:CA	2.91	0.49
1:B:577:LEU:HD23	1:B:687:ILE:HD13	1.95	0.49
1:A:796:VAL:HB	1:A:837:LEU:HD22	1.94	0.48
1:A:556:GLN:HE22	1:B:579:SER:HA	1.78	0.48
1:A:795:LEU:HD23	1:A:836:VAL:HB	1.97	0.47
1:B:697:LYS:O	1:B:698:ALA:HB3	2.15	0.46
1:C:792:ASP:O	1:C:831:ALA:HB1	2.16	0.46
1:C:757:VAL:HG22	1:C:891:HIS:CE1	2.50	0.46
1:A:810:LEU:HD22	4:A:1900:NU7:C29	2.46	0.46
1:A:824:THR:CG2	1:A:859:LEU:HD13	2.46	0.45
1:A:520:THR:HB	1:A:659:ASN:OD1	2.16	0.45
1:C:824:THR:HG21	1:C:859:LEU:HD13	1.99	0.44
1:A:525:ASP:OD1	1:A:527:VAL:HG23	2.17	0.44
1:A:577:LEU:HD23	1:A:687:ILE:HD13	1.99	0.44
1:C:628:ILE:HG22	1:C:629:TRP:N	2.31	0.44
1:B:702:LEU:HD13	1:B:725:LEU:HD23	2.00	0.44
1:B:810:LEU:HD22	4:B:1900:NU7:C29	2.47	0.44
1:A:523:ILE:HG21	1:A:644:SER:HB3	1.99	0.44
1:A:817:ALA:HB2	1:A:850:ASP:HB3	2.00	0.43
1:B:548:ILE:HD13	1:B:665:ARG:HG2	2.01	0.43
1:C:592:ASN:HB3	1:C:593:PRO:HA	2.00	0.43
1:C:560:LEU:HD13	1:C:856:VAL:HG12	2.01	0.43
1:A:649:ALA:HA	1:A:836:VAL:HG21	2.01	0.43
1:B:731:ARG:CZ	1:B:765:LEU:HD11	2.49	0.42
1:C:554:ARG:NH2	1:C:850:ASP:OD1	2.53	0.42
1:C:529:LEU:HD23	1:C:548:ILE:HD11	2.01	0.41
1:B:592:ASN:O	1:B:595:SER:N	2.53	0.41
1:A:824:THR:HG21	1:A:859:LEU:HD13	2.02	0.41
1:C:592:ASN:HB3	1:C:593:PRO:CA	2.49	0.41
1:B:796:VAL:CB	1:B:837:LEU:HD22	2.51	0.40
1:A:757:VAL:HG22	1:A:891:HIS:CE1	2.56	0.40
1:B:796:VAL:CG2	1:B:837:LEU:HD22	2.51	0.40
1:C:824:THR:CG2	1:C:859:LEU:HD13	2.52	0.40
1:B:757:VAL:HG22	1:B:891:HIS:CE1	2.56	0.40
1:C:541:HIS:ND1	1:C:628:ILE:HD11	2.37	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	364/423 (86%)	347 (95%)	16 (4%)	1 (0%)	46	57
1	B	376/423 (89%)	363 (96%)	12 (3%)	1 (0%)	46	57
1	C	355/423 (84%)	341 (96%)	11 (3%)	3 (1%)	24	28
All	All	1095/1269 (86%)	1051 (96%)	39 (4%)	5 (0%)	34	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	592	ASN
1	A	811	GLY
1	C	811	GLY
1	B	811	GLY
1	C	673	HIS

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	289/336 (86%)	277 (96%)	12 (4%)	36	51
1	B	298/336 (89%)	284 (95%)	14 (5%)	32	45
1	C	280/336 (83%)	276 (99%)	4 (1%)	74	85
All	All	867/1008 (86%)	837 (96%)	30 (4%)	43	60

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	519	THR
1	A	520	THR
1	A	532	GLN
1	A	543	GLU
1	A	553	SER
1	A	647	ASP
1	A	657	LEU
1	A	669	HIS
1	A	692	LEU
1	A	706	TRP
1	A	734	ASP
1	A	859	LEU
1	B	543	GLU
1	B	561	ARG
1	B	574	LEU
1	B	594	LEU
1	B	623	VAL
1	B	647	ASP
1	B	669	HIS
1	B	706	TRP
1	B	765	LEU
1	B	766	ASP
1	B	839	LEU
1	B	862	ASN
1	B	874	GLN
1	B	878	LEU
1	C	523	ILE
1	C	561	ARG
1	C	669	HIS
1	C	765	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	556	GLN
1	A	636	ASN
1	A	673	HIS
1	A	693	GLN
1	A	720	GLN
1	A	756	ASN
1	A	879	ASN
1	A	891	HIS
1	A	899	GLN

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Mol	Chain	Res	Type
1	B	556	GLN
1	B	636	ASN
1	B	825	GLN
1	B	879	ASN
1	B	899	GLN
1	C	636	ASN
1	C	673	HIS
1	C	693	GLN
1	C	720	GLN
1	C	899	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 12 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	NU7	A	1900	-	35,41,41	1.14	2 (5%)	39,60,60	2.02	7 (17%)
4	NU7	B	1900	-	35,41,41	1.11	2 (5%)	39,60,60	2.10	6 (15%)
4	NU7	C	1900	2	35,41,41	1.12	2 (5%)	39,60,60	2.13	6 (15%)

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	NU7	A	1900	-	-	0/21/42/42	0/4/5/5
4	NU7	B	1900	-	-	0/21/42/42	0/4/5/5
4	NU7	C	1900	2	-	0/21/42/42	0/4/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1900	NU7	C07-N06	3.07	1.39	1.35
4	A	1900	NU7	C07-N06	3.38	1.40	1.35
4	B	1900	NU7	C07-N06	3.46	1.40	1.35
4	B	1900	NU7	C29-S30	3.94	1.76	1.70
4	C	1900	NU7	C29-S30	4.29	1.77	1.70
4	A	1900	NU7	C29-S30	4.29	1.77	1.70

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1900	NU7	C29-C28-C31	-6.17	120.81	129.40
4	C	1900	NU7	C29-C28-C31	-5.28	122.05	129.40
4	B	1900	NU7	C28-C29-S30	-5.26	105.33	111.79
4	A	1900	NU7	C29-C28-C31	-5.21	122.14	129.40
4	C	1900	NU7	C28-C29-S30	-5.02	105.62	111.79
4	A	1900	NU7	C28-C29-S30	-4.99	105.66	111.79
4	C	1900	NU7	C20-C19-N18	-2.50	109.17	113.88
4	A	1900	NU7	C20-C19-N18	-2.09	109.94	113.88
4	A	1900	NU7	C21-C22-N2	2.02	112.50	110.88
4	B	1900	NU7	C21-C22-N2	2.48	112.86	110.88
4	C	1900	NU7	C21-C22-N2	2.88	113.18	110.88
4	B	1900	NU7	C02-C05-N06	3.51	125.83	122.41
4	A	1900	NU7	C02-C05-N06	4.43	126.73	122.41
4	A	1900	NU7	C10-C07-N08	5.00	125.41	119.11
4	B	1900	NU7	C10-C07-N08	5.33	125.83	119.11
4	A	1900	NU7	C05-N06-C07	5.70	108.12	101.45
4	B	1900	NU7	C05-N06-C07	6.01	108.49	101.45
4	C	1900	NU7	C05-N06-C07	6.31	108.84	101.45
4	C	1900	NU7	C10-C07-N08	6.90	127.81	119.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1900	NU7	1	0
4	B	1900	NU7	2	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

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	368/423 (86%)	0.61	44 (11%) 6 5	23, 29, 35, 39	0
1	B	380/423 (89%)	0.78	61 (16%) 3 2	23, 29, 35, 45	0
1	C	359/423 (84%)	0.91	62 (17%) 2 2	28, 33, 38, 43	0
All	All	1107/1269 (87%)	0.77	167 (15%) 3 3	23, 31, 36, 45	0

All (167) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	796	VAL	6.8
1	B	596	ARG	6.4
1	B	594	LEU	5.9
1	B	704	VAL	5.9
1	C	798	ALA	5.5
1	B	797	SER	5.3
1	A	796	VAL	5.2
1	C	562	SER	5.1
1	C	704	VAL	5.1
1	B	706	TRP	5.0
1	A	706	TRP	4.9
1	B	796	VAL	4.9
1	B	685	VAL	4.9
1	B	798	ALA	4.7
1	C	837	LEU	4.6
1	C	867	LEU	4.6
1	C	838	ALA	4.5
1	C	564	CYS	4.4
1	A	704	VAL	4.4
1	C	797	SER	4.3
1	A	838	ALA	4.3
1	C	799	GLY	4.1
1	B	799	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	838	ALA	4.1
1	A	837	LEU	4.1
1	C	655	ARG	4.0
1	A	797	SER	3.9
1	C	809	PRO	3.9
1	C	616	LEU	3.9
1	A	798	ALA	3.8
1	A	809	PRO	3.8
1	B	537	ASP	3.8
1	B	593	PRO	3.7
1	C	703	ILE	3.7
1	B	539	SER	3.7
1	B	705	ASP	3.7
1	B	729	LEU	3.7
1	C	657	LEU	3.6
1	C	539	SER	3.6
1	A	729	LEU	3.6
1	C	808	ALA	3.5
1	C	560	LEU	3.5
1	B	668	GLY	3.5
1	B	839	LEU	3.5
1	B	610	GLN	3.3
1	C	705	ASP	3.3
1	B	669	HIS	3.3
1	C	653	ALA	3.3
1	A	664	VAL	3.2
1	B	711	GLY	3.2
1	C	706	TRP	3.2
1	C	752	GLY	3.1
1	A	685	VAL	3.1
1	B	597	LEU	3.1
1	A	778	PHE	3.1
1	B	664	VAL	3.1
1	C	871	GLY	3.0
1	B	592	ASN	3.0
1	C	805	GLY	3.0
1	C	795	LEU	3.0
1	A	705	ASP	3.0
1	A	827	LEU	2.9
1	B	727	ILE	2.9
1	B	863	ARG	2.9
1	B	837	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	841	GLY	2.9
1	B	728	SER	2.9
1	A	662	ALA	2.9
1	C	568	ARG	2.9
1	A	684	SER	2.9
1	C	860	LEU	2.9
1	A	666	PRO	2.9
1	A	703	ILE	2.9
1	C	841	GLY	2.9
1	C	685	VAL	2.8
1	A	836	VAL	2.8
1	B	734	ASP	2.8
1	C	734	ASP	2.8
1	B	680	CYS	2.8
1	B	823	MET	2.8
1	B	666	PRO	2.8
1	C	617	PRO	2.7
1	A	823	MET	2.7
1	A	645	VAL	2.7
1	B	795	LEU	2.7
1	C	810	LEU	2.7
1	B	800	PHE	2.6
1	B	870	GLU	2.6
1	B	684	SER	2.6
1	A	765	LEU	2.6
1	B	518	PHE	2.6
1	B	774	TYR	2.6
1	A	736	ASN	2.6
1	A	702	LEU	2.5
1	C	836	VAL	2.5
1	B	667	PRO	2.5
1	C	768	PRO	2.5
1	C	656	GLU	2.5
1	C	557	GLU	2.5
1	A	516	LEU	2.5
1	A	795	LEU	2.5
1	C	549	GLN	2.5
1	C	711	GLY	2.5
1	A	686	ALA	2.5
1	C	699	SER	2.5
1	B	809	PRO	2.4
1	B	782	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	790	SER	2.4
1	A	616	LEU	2.4
1	A	839	LEU	2.4
1	C	727	ILE	2.4
1	A	808	ALA	2.4
1	A	766	ASP	2.4
1	B	617	PRO	2.4
1	C	558	ARG	2.4
1	A	824	THR	2.4
1	C	751	SER	2.4
1	C	647	ASP	2.4
1	A	800	PHE	2.4
1	A	840	GLU	2.4
1	A	868	SER	2.4
1	B	703	ILE	2.4
1	B	726	TYR	2.4
1	C	818	LYS	2.4
1	C	874	GLN	2.4
1	C	840	GLU	2.4
1	B	536	GLY	2.3
1	B	595	SER	2.3
1	C	519	THR	2.3
1	B	777	ALA	2.3
1	A	660	GLY	2.3
1	B	780	ILE	2.3
1	C	658	LYS	2.3
1	C	873	LYS	2.3
1	A	820	PHE	2.3
1	C	521	GLY	2.3
1	A	641	ALA	2.2
1	C	537	ASP	2.2
1	B	707	ASP	2.2
1	C	623	VAL	2.2
1	C	668	GLY	2.2
1	C	839	LEU	2.2
1	B	540	ARG	2.1
1	B	714	THR	2.1
1	B	557	GLU	2.1
1	B	654	SER	2.1
1	C	553	SER	2.1
1	B	653	ALA	2.1
1	B	865	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	667	PRO	2.1
1	A	595	SER	2.1
1	C	662	ALA	2.1
1	A	683	ASN	2.1
1	A	841	GLY	2.1
1	B	620	GLY	2.1
1	C	625	THR	2.1
1	C	870	GLU	2.1
1	B	867	LEU	2.1
1	C	722	PRO	2.1
1	B	840	GLU	2.0
1	B	702	LEU	2.0
1	C	736	ASN	2.0
1	A	661	PHE	2.0
1	B	846	THR	2.0
1	B	768	PRO	2.0
1	A	725	LEU	2.0
1	C	592	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	K	C	202	1/1	0.70	0.23	1.13	86,86,86,86	0
3	K	B	202	1/1	0.89	0.22	0.82	54,54,54,54	0
3	K	C	201	1/1	0.92	0.30	0.82	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NU7	B	1900	37/37	0.71	0.30	0.79	64,71,74,74	0
4	NU7	C	1900	37/37	0.75	0.28	0.69	83,85,86,86	0
3	K	A	202	1/1	0.82	0.19	0.45	65,65,65,65	0
3	K	B	201	1/1	0.95	0.34	0.43	61,61,61,61	0
4	NU7	A	1900	37/37	0.80	0.22	-0.01	64,65,68,68	0
3	K	A	201	1/1	0.93	0.17	-1.23	61,61,61,61	0
2	ZN	B	101	1/1	0.99	0.22	-1.26	34,34,34,34	0
2	ZN	B	102	1/1	0.98	0.03	-2.05	24,24,24,24	0
2	ZN	C	102	1/1	0.89	0.08	-2.16	33,33,33,33	0
2	ZN	C	101	1/1	0.98	0.11	-2.36	32,32,32,32	0
2	ZN	A	102	1/1	0.97	0.03	-2.66	30,30,30,30	0
2	ZN	A	101	1/1	0.98	0.14	-3.28	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.