Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:34 PM GMT

PDB ID	:	3U3Y
Title	:	Mouse TREX1 D200H mutant
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Deposited on	:	2011-10-06
Resolution	:	2.28 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 (i) symbol.

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

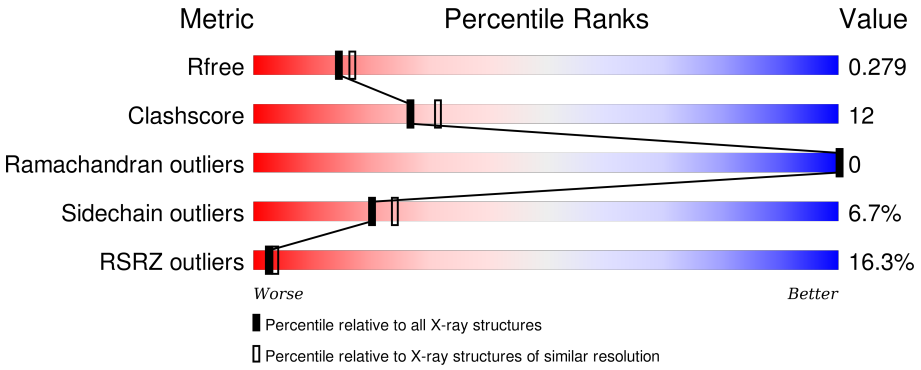
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.28 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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 <=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	314	11% 52% 14% • 32%
1	B	314	12% 52% 16% • 30%
2	C	4	25% 100%
2	D	4	50% 50%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BU1	A	315	-	-	X	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7123 atoms, of which 3388 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 Three prime repair exonuclease 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	220	Total	C	H	N	O	S	0	1	0
			3408	1082	1702	301	314	9			
1	A	215	Total	C	H	N	O	S	0	1	0
			3352	1066	1676	295	306	9			

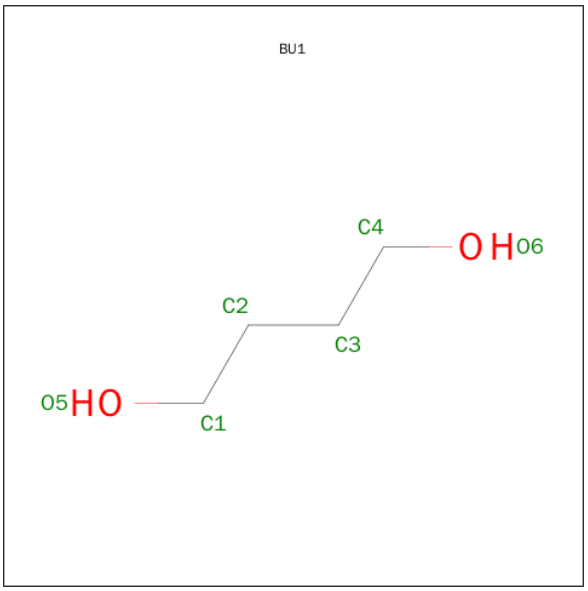
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	200	HIS	ASP	ENGINEERED MUTATION	UNP Q91XB0
A	200	HIS	ASP	ENGINEERED MUTATION	UNP Q91XB0

- Molecule 2 is a DNA chain called 5'-D(*GP*AP*CP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	4	Total	C	N	O	P	0	0	0
			81	39	18	21	3			
2	D	4	Total	C	N	O	P	0	0	0
			81	39	18	21	3			

- Molecule 3 is 1,4-BUTANEDIOL (three-letter code: BU1) (formula: $C_4H_{10}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 16 4 10 2	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	Total Ca 1 1	0	0
4	C	1	Total Ca 1 1	0	0

- Molecule 5 is water.

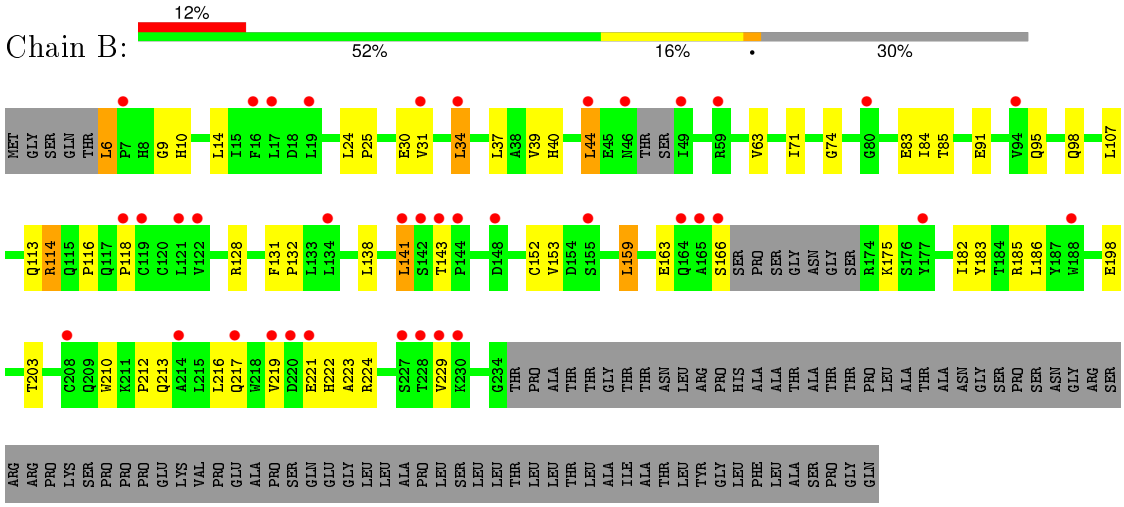
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	76	Total O 76 76	0	0
5	A	88	Total O 88 88	0	0
5	C	7	Total O 7 7	0	0
5	D	12	Total O 12 12	0	0

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 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.

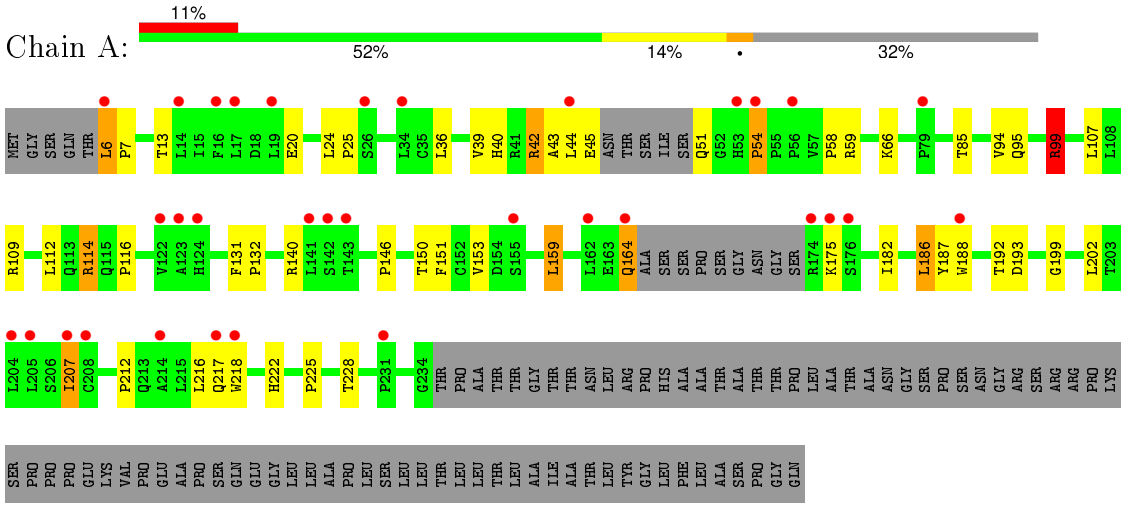
- Molecule 1: Three prime repair exonuclease 1

Chain B:



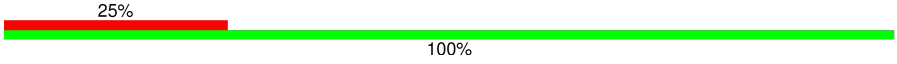
- Molecule 1: Three prime repair exonuclease 1

Chain A:



- Molecule 2: 5'-D(*GP*AP*CP*G)-3'

Chain C:



- Molecule 2: 5'-D(*GP*AP*CP*G)-3'

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	65.07Å 57.86Å 68.23Å 90.00° 108.11° 90.00°	Depositor
Resolution (Å)	22.72 – 2.28 22.72 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.2 (22.72-2.28) 99.4 (22.72-2.28)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.42 (at 2.28Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.213 , 0.275 0.221 , 0.279	Depositor DCC
R_{free} test set	1137 reflections (5.43%)	DCC
Wilson B-factor (Å^2)	33.7	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 51.0	EDS
Estimated twinning fraction	0.012 for l,-k,h	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 22059 reflections	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7123	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% 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, 0.375 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: CA, BU1

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	0.40	1/1721 (0.1%)	0.58	1/2345 (0.0%)
1	B	0.36	0/1751	0.54	0/2386
2	C	0.74	0/91	1.14	0/139
2	D	0.74	0/91	1.20	0/139
All	All	0.40	1/3654 (0.0%)	0.61	1/5009 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	54	PRO	C-N	-5.47	1.23	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	164	GLN	N-CA-CB	-5.52	100.67	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	99	ARG	Sidechain

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	1676	1676	1678	47	0
1	B	1706	1702	1701	45	0
2	C	81	0	46	0	0
2	D	81	0	46	2	0
3	A	6	10	10	10	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	88	0	0	1	0
5	B	76	0	0	1	0
5	C	7	0	0	0	0
5	D	12	0	0	1	0
All	All	3735	3388	3481	85	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.

All (85) 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:A:202:LEU:HD12	3:A:315:BU1:H22	1.57	0.86
1:A:212:PRO:O	1:A:216:LEU:HD23	1.89	0.72
1:A:159:LEU:HD22	1:A:182:ILE:HG21	1.71	0.72
1:A:54:PRO:HD3	1:A:188:TRP:CZ3	2.26	0.71
1:A:192:THR:HB	3:A:315:BU1:C3	2.22	0.70
1:B:159:LEU:HD22	1:B:182:ILE:HG21	1.73	0.70
1:A:164:GLN:OE1	1:A:164:GLN:N	2.24	0.69
1:A:54:PRO:HD3	1:A:188:TRP:CH2	2.32	0.65
1:B:198:GLU:OE1	1:A:66:LYS:NZ	2.28	0.64
1:B:44:LEU:O	1:B:212:PRO:HB3	1.99	0.62
1:B:185:ARG:HH11	1:B:185:ARG:HG3	1.66	0.60
1:B:221:GLU:HG3	1:B:222:HIS:ND1	2.16	0.60
1:A:6:LEU:HB2	1:A:7:PRO:HD3	1.85	0.59
1:A:202:LEU:HB2	3:A:315:BU1:H42	1.84	0.58
1:A:109:ARG:HB2	1:A:146:PRO:HB3	1.85	0.58
1:A:193:ASP:HB3	3:A:315:BU1:H12	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:LEU:HD12	1:A:107:LEU:CD1	2.33	0.58
1:B:40:HIS:HD2	1:B:116:PRO:HG2	1.70	0.56
1:B:107:LEU:CD1	1:A:107:LEU:CD1	2.83	0.56
1:B:37:LEU:CD2	1:B:39:VAL:HG23	2.35	0.56
1:B:175:LYS:HD2	1:B:185:ARG:HD3	1.87	0.56
1:A:94:VAL:HG12	1:A:95:GLN:HG3	1.87	0.56
1:A:131:PHE:HB2	1:A:132:PRO:HD3	1.88	0.56
1:B:131:PHE:HB2	1:B:132:PRO:HD3	1.88	0.55
1:A:140:ARG:HD2	5:A:337:HOH:O	2.06	0.55
1:A:58:PRO:HD3	1:A:187:TYR:CE2	2.43	0.54
1:A:182:ILE:HG22	1:A:186:LEU:HD22	1.90	0.53
1:A:192:THR:HB	3:A:315:BU1:C4	2.39	0.53
1:B:128:ARG:NH1	2:D:1:DG:N7	2.57	0.53
1:B:114:ARG:NH2	1:A:99:ARG:O	2.41	0.53
1:A:202:LEU:CD1	3:A:315:BU1:H22	2.35	0.52
1:B:24:LEU:HB3	1:B:25:PRO:HD2	1.91	0.52
1:B:183:TYR:CD2	1:B:203:THR:HG23	2.44	0.52
1:A:212:PRO:O	1:A:216:LEU:CD2	2.58	0.51
1:B:44:LEU:O	1:B:212:PRO:CB	2.59	0.50
1:B:131:PHE:CE1	1:B:152:CYS:HB2	2.46	0.50
1:B:91:GLU:HG3	5:B:377:HOH:O	2.12	0.50
1:B:95:GLN:OE1	1:A:43:ALA:HB2	2.12	0.50
1:A:40:HIS:CE1	1:A:42:ARG:HG2	2.47	0.49
1:A:159:LEU:HD11	1:A:207:ILE:CG2	2.43	0.49
1:B:63:VAL:HB	1:A:95:GLN:OE1	2.13	0.48
1:B:107:LEU:CD1	1:A:107:LEU:HD12	2.44	0.48
1:B:153:VAL:CG1	1:B:223:ALA:HA	2.44	0.48
1:A:225:PRO:HB2	1:A:228:THR:HG23	1.95	0.48
1:B:221:GLU:HG3	1:B:222:HIS:HD1	1.78	0.48
1:B:37:LEU:HD23	1:B:37:LEU:C	2.35	0.47
1:A:42:ARG:HA	1:A:45:GLU:HG2	1.95	0.47
1:B:224:ARG:HH21	1:B:229:VAL:HG22	1.79	0.47
1:B:34:LEU:HD23	1:B:34:LEU:C	2.34	0.47
1:A:39:VAL:HG12	1:A:43:ALA:HB3	1.97	0.47
1:A:54:PRO:HD3	1:A:188:TRP:CE3	2.50	0.46
1:A:199:GLY:HA2	3:A:315:BU1:H12	1.97	0.46
1:B:40:HIS:CD2	1:B:116:PRO:HG2	2.52	0.45
1:B:98:GLN:O	1:A:114:ARG:NH2	2.48	0.45
1:B:131:PHE:CZ	1:B:152:CYS:HB2	2.52	0.44
1:A:192:THR:HB	3:A:315:BU1:H32	1.96	0.44
1:B:14:LEU:HD11	1:B:216:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:GLU:OE1	1:B:74:GLY:N	2.46	0.43
1:B:186:LEU:HD13	1:B:210:TRP:CG	2.54	0.43
1:A:40:HIS:ND1	1:A:42:ARG:HG2	2.33	0.43
1:A:159:LEU:HD11	1:A:207:ILE:HG23	2.00	0.43
1:B:182:ILE:O	1:B:186:LEU:HG	2.19	0.43
1:A:159:LEU:CD1	1:A:207:ILE:CG2	2.97	0.43
1:B:24:LEU:HB3	1:B:25:PRO:CD	2.49	0.43
1:B:138:LEU:O	1:B:141:LEU:HB2	2.19	0.43
1:B:163:GLU:O	1:B:166:SER:N	2.48	0.43
1:A:193:ASP:HB3	3:A:315:BU1:H31	2.01	0.42
1:A:20[B]:GLU:OE1	1:A:85:THR:HB	2.18	0.42
1:A:218:TRP:CD1	1:A:222:HIS:CE1	3.07	0.42
1:B:84:ILE:HG22	1:B:85:THR:HG23	2.02	0.42
1:B:185:ARG:NH1	1:B:185:ARG:HG3	2.34	0.41
1:A:44:LEU:CB	1:A:216:LEU:HD21	2.50	0.41
1:B:153:VAL:HG13	1:B:223:ALA:HA	2.03	0.41
1:B:6:LEU:HD23	1:B:9:GLY:C	2.40	0.41
1:B:6:LEU:HG	1:B:9:GLY:O	2.21	0.41
1:B:10:HIS:O	1:B:118:PRO:HD2	2.21	0.41
1:A:193:ASP:CB	3:A:315:BU1:H31	2.51	0.41
1:B:153:VAL:HG21	1:B:219:VAL:HG22	2.02	0.41
1:A:150:THR:HG22	1:A:151:PHE:N	2.35	0.41
1:A:24:LEU:HB3	1:A:25:PRO:HD2	2.03	0.41
1:A:164:GLN:CD	1:A:164:GLN:N	2.74	0.40
2:D:3:DC:H5	5:D:177:HOH:O	2.00	0.40
1:B:31:VAL:HG12	1:B:71:ILE:HD12	2.04	0.40
1:B:14:LEU:HD11	1:B:216:LEU:CD2	2.52	0.40
1:A:13:THR:OG1	1:A:116:PRO:HD2	2.22	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	210/314 (67%)	205 (98%)	5 (2%)	0	100	100
1	B	215/314 (68%)	213 (99%)	2 (1%)	0	100	100
All	All	425/628 (68%)	418 (98%)	7 (2%)	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	185/263 (70%)	171 (92%)	14 (8%)	16	19
1	B	188/263 (72%)	177 (94%)	11 (6%)	24	30
All	All	373/526 (71%)	348 (93%)	25 (7%)	20	24

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	6	LEU
1	B	34	LEU
1	B	44	LEU
1	B	83	GLU
1	B	113	GLN
1	B	114	ARG
1	B	141	LEU
1	B	143	THR
1	B	159	LEU
1	B	213	GLN
1	B	217	GLN
1	A	6	LEU
1	A	36	LEU
1	A	42	ARG
1	A	51	GLN
1	A	59	ARG
1	A	99	ARG
1	A	112	LEU

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Mol	Chain	Res	Type
1	A	114	ARG
1	A	153	VAL
1	A	159	LEU
1	A	175	LYS
1	A	186	LEU
1	A	207	ILE
1	A	217	GLN

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 molecules 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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	BU1	A	315	-	5,5,5	0.33	0	4,4,4	0.36	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
3	BU1	A	315	-	-	0/3/3/3	0/0/0/0

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.

1 monomer is involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	315	BU1	10	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, 95^{th} 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	215/314 (68%)	0.89	33 (15%) 3 4	20, 38, 66, 99	0
1	B	220/314 (70%)	0.93	38 (17%) 2 3	19, 41, 72, 88	0
2	C	4/4 (100%)	0.51	1 (25%) 1 1	28, 36, 43, 49	0
2	D	4/4 (100%)	0.23	0 100 100	36, 37, 37, 50	0
All	All	443/636 (69%)	0.90	72 (16%) 2 4	19, 40, 69, 99	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	165	ALA	6.2
1	B	164	GLN	5.9
1	A	188	TRP	5.3
1	A	142	SER	5.1
1	A	208	CYS	4.8
1	B	142	SER	4.4
1	B	94	VAL	4.4
1	A	204	LEU	4.3
1	A	205	LEU	4.2
1	B	208	CYS	4.2
1	A	217	GLN	4.1
1	A	17	LEU	3.9
1	A	79	PRO	3.9
1	A	155	SER	3.6
1	B	49	ILE	3.6
1	B	228	THR	3.5
1	A	16	PHE	3.4
1	A	175	LYS	3.4
1	A	54	PRO	3.4
1	A	207	ILE	3.4
1	A	6	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	143	THR	3.3
1	A	174	ARG	3.2
1	B	122	VAL	3.2
1	A	122	VAL	3.2
1	A	19	LEU	3.1
1	B	229	VAL	3.1
1	B	227	SER	3.0
1	B	148	ASP	2.9
1	A	162	LEU	2.9
1	B	134	LEU	2.9
1	B	141	LEU	2.9
1	B	166	SER	2.9
1	B	144	PRO	2.8
1	B	59	ARG	2.8
1	B	16	PHE	2.8
1	B	19	LEU	2.8
1	A	124	HIS	2.7
1	A	123	ALA	2.6
1	A	53	HIS	2.6
1	A	56	PRO	2.6
1	B	44	LEU	2.6
1	A	14	LEU	2.6
1	B	80	GLY	2.5
1	B	230	LYS	2.5
1	B	119	CYS	2.5
1	B	46	ASN	2.5
1	B	155	SER	2.5
1	B	220	ASP	2.5
1	B	214	ALA	2.4
1	A	141	LEU	2.4
2	C	1	DG	2.4
1	A	176	SER	2.4
1	A	143	THR	2.4
1	B	34	LEU	2.4
1	B	217	GLN	2.4
1	B	118	PRO	2.4
1	B	17	LEU	2.3
1	B	219	VAL	2.3
1	A	34	LEU	2.3
1	A	44	LEU	2.3
1	A	218	TRP	2.3
1	B	7	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	31	VAL	2.2
1	B	177	TYR	2.2
1	A	231	PRO	2.2
1	A	164	GLN	2.1
1	A	26	SER	2.1
1	B	121	LEU	2.1
1	A	214	ALA	2.1
1	B	188	TRP	2.0
1	B	221	GLU	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, 95^{th} 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	BU1	A	315	6/6	0.84	0.41	6.58	30,39,42,50	0
4	CA	C	5	1/1	0.98	0.10	-1.75	47,47,47,47	0
4	CA	D	5	1/1	0.98	0.04	-3.28	35,35,35,35	0

6.5 Other polymers (i)

There are no such residues in this entry.