Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:35 PM GMT

PDB ID	:	1BIQ
Title	:	RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 BETA CHAIN MUTANT E238A
Authors	:	Logan, D.T.; Demare, F.; Persson, B.O.; Slaby, A.; Sjoberg, B.M.; Nordlund, P.
Deposited on	:	1998-06-18
Resolution	:	2.05 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:

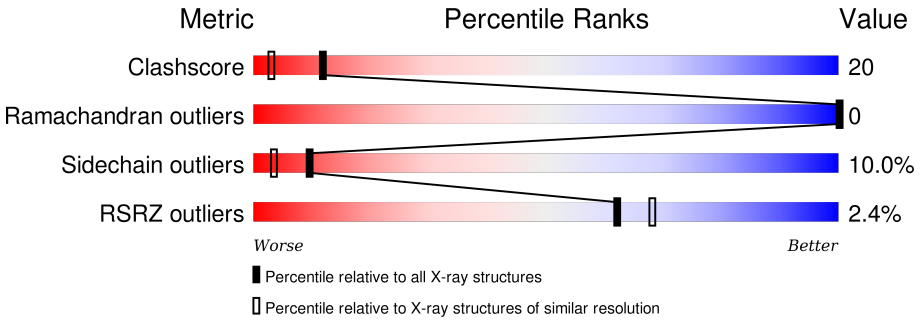
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

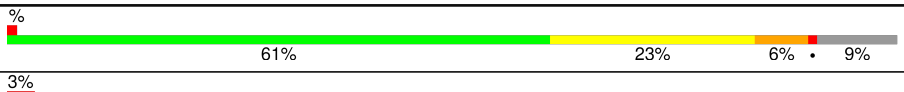
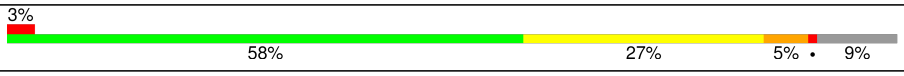
The reported resolution of this entry is 2.05 Å.

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))
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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	375	
2	B	375	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5921 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 PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	1	0
			2784	1782	464	525	13			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	122	PHE	TYR	ENGINEERED	UNP P69924
A	208	MTY	PHE	MODIFIED RESIDUE	UNP P69924
A	238	ALA	GLU	ENGINEERED	UNP P69924

- Molecule 2 is a protein called PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	341	Total	C	N	O	S	0	2	0
			2789	1785	465	526	13			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	122	PHE	TYR	ENGINEERED	UNP P69924
B	238	ALA	GLU	ENGINEERED	UNP P69924

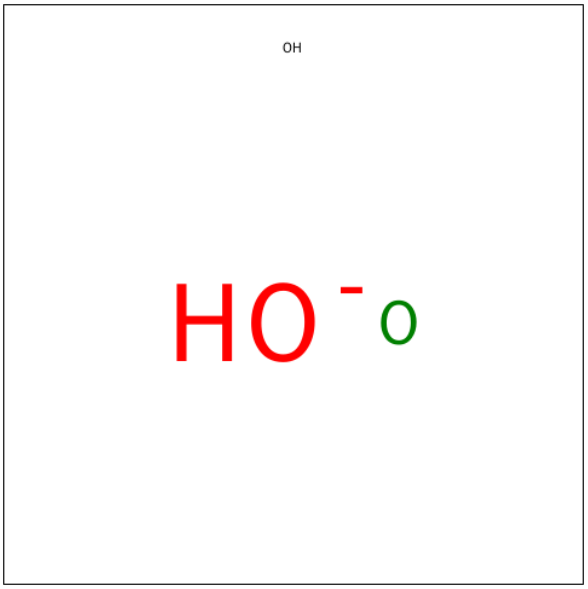
- Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Fe	0	0
			2	2		
3	A	1	Total	Fe	0	0
			1	1		

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Fe 1 1	0	0

- Molecule 5 is HYDROXIDE ION (three-letter code: OH) (formula: HO).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O 1 1	0	0
5	A	1	Total O 1 1	0	0
5	B	1	Total O 1 1	0	0

- Molecule 6 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	9	Total Hg 9 9	0	0
6	A	7	Total Hg 7 7	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	160	Total O 160 160	0	0

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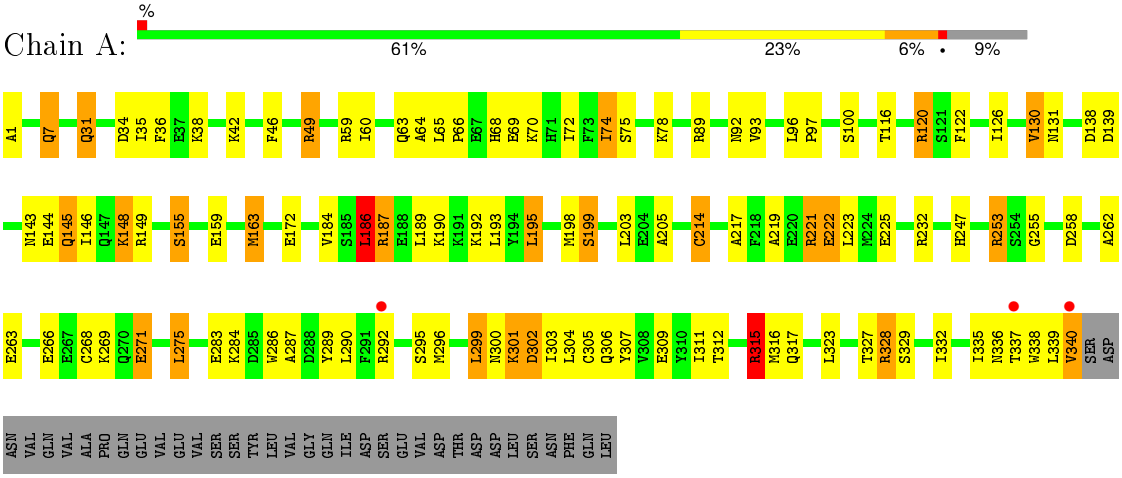
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	165	Total	O	0	0
			165	165		

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.

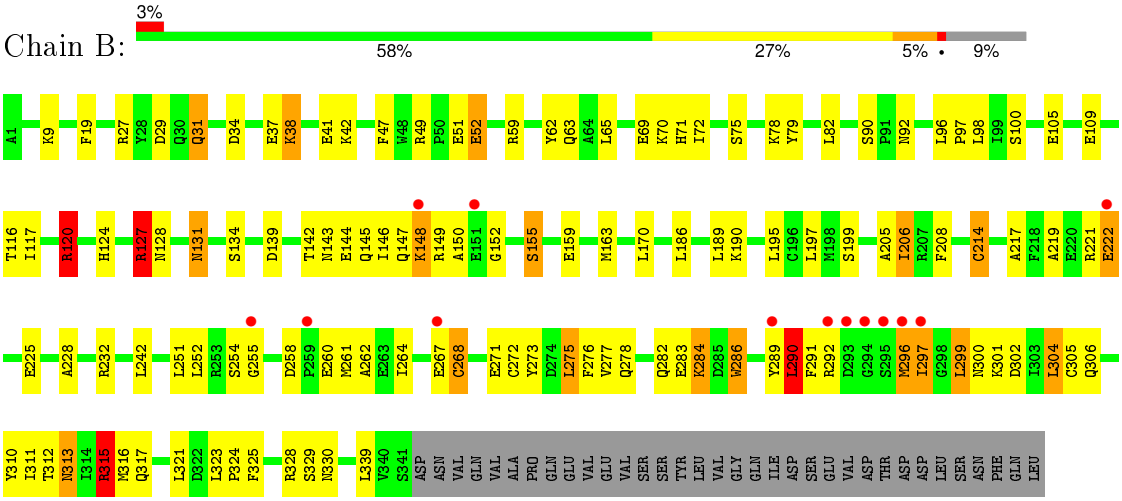
- Molecule 1: PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE

Chain A:



- Molecule 2: PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	74.05Å 83.78Å 113.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.05 27.12 – 2.06	Depositor EDS
% Data completeness (in resolution range)	98.0 (25.00-2.05) 98.2 (27.12-2.06)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	13.32 (at 2.06Å)	Xtriage
Refinement program	TNT V. 5-E	Depositor
R, R_{free}	0.191 , 0.260 0.182 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	22.0	Xtriage
Anisotropy	0.456	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 98.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 43712 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5921	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 5.52% 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: HG, MTY, FE2, FE, OH

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	0.80	1/2834 (0.0%)	1.23	18/3843 (0.5%)
2	B	0.75	0/2853	1.18	9/3870 (0.2%)
All	All	0.77	1/5687 (0.0%)	1.20	27/7713 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	269	LYS	CE-NZ	6.96	1.66	1.49

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	315	ARG	NE-CZ-NH1	13.58	127.09	120.30
1	A	340	VAL	CG1-CB-CG2	8.77	124.93	110.90
1	A	89	ARG	NE-CZ-NH1	7.04	123.82	120.30
2	B	120	ARG	NE-CZ-NH1	-6.74	116.93	120.30
1	A	120	ARG	NE-CZ-NH1	6.63	123.61	120.30
2	B	290	LEU	CB-CG-CD2	-6.51	99.93	111.00
2	B	127	ARG	NE-CZ-NH1	-6.43	117.09	120.30
1	A	214	CYS	CB-CA-C	6.24	122.89	110.40
1	A	275	LEU	CB-CG-CD2	-6.12	100.59	111.00
1	A	187	ARG	NE-CZ-NH1	-6.08	117.26	120.30
1	A	198	MET	CG-SD-CE	6.06	109.89	100.20
1	A	315	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	A	328	ARG	NE-CZ-NH2	-5.95	117.32	120.30
2	B	251	LEU	CB-CG-CD1	5.91	121.04	111.00
1	A	275	LEU	CB-CG-CD1	-5.62	101.44	111.00
2	B	339	LEU	CB-CG-CD2	-5.42	101.78	111.00
2	B	90	SER	N-CA-C	5.29	125.29	111.00
1	A	221	ARG	NE-CZ-NH1	-5.27	117.67	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	195	LEU	CB-CG-CD2	5.23	119.90	111.00
1	A	193	LEU	CB-CG-CD1	5.23	119.89	111.00
1	A	163	MET	CG-SD-CE	5.20	108.52	100.20
2	B	299	LEU	CA-CB-CG	5.13	127.09	115.30
2	B	315	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	A	232	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	A	328	ARG	NE-CZ-NH1	5.06	122.83	120.30
2	B	214[A]	CYS	CA-CB-SG	-5.04	104.93	114.00
1	A	186	LEU	CB-CG-CD2	5.02	119.53	111.00

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	2784	0	2729	97	0
2	B	2789	0	2732	121	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	1	0	0	0	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0
6	A	7	0	0	1	0
6	B	9	0	0	2	0
7	A	160	0	0	9	0
7	B	165	0	0	13	0
All	All	5921	0	5461	215	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 20.

All (215) 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:205:ALA:HB1	1:A:315:ARG:HD3	1.26	1.17
1:A:138:ASP:HB3	2:B:9:LYS:HE2	1.21	1.07
2:B:72:ILE:HG23	2:B:290:LEU:HD21	1.53	0.91
1:A:332:ILE:HD12	1:A:335:ILE:HG21	1.52	0.91
2:B:31:GLN:HG3	2:B:34:ASP:HA	1.53	0.88
1:A:287:ALA:CB	1:A:301:LYS:HD3	2.05	0.86
1:A:287:ALA:HB3	1:A:301:LYS:HD3	1.55	0.86
2:B:214[A]:CYS:SG	6:B:387:HG:HG	1.92	0.86
1:A:301:LYS:HD2	1:A:305[A]:CYS:SG	2.16	0.85
1:A:205:ALA:HB1	1:A:315:ARG:CD	2.07	0.83
2:B:149:ARG:HD2	2:B:286:TRP:CG	2.13	0.83
2:B:272[A]:CYS:SG	6:B:380:HG:HG	1.97	0.83
2:B:72:ILE:HG12	2:B:290:LEU:HD23	1.62	0.82
2:B:149:ARG:HD2	2:B:286:TRP:CD2	2.16	0.80
1:A:42:LYS:HE2	1:A:46:PHE:CZ	2.16	0.80
1:A:332:ILE:HD12	1:A:335:ILE:CG2	2.12	0.80
1:A:205:ALA:CB	1:A:315:ARG:HD3	2.12	0.77
1:A:138:ASP:HB3	2:B:9:LYS:CE	2.09	0.76
2:B:205:ALA:HB2	2:B:242:LEU:CD2	2.15	0.75
2:B:163:MET:HE2	2:B:189:LEU:HB2	1.68	0.75
2:B:221:ARG:CZ	2:B:297:ILE:HG13	2.17	0.74
1:A:42:LYS:HE2	1:A:46:PHE:HZ	1.51	0.74
2:B:252:LEU:HD22	2:B:261:MET:HG2	1.69	0.73
2:B:271:GLU:O	2:B:275:LEU:HD22	1.87	0.73
2:B:205:ALA:HB2	2:B:242:LEU:HD23	1.70	0.73
7:A:544:HOH:O	2:B:49:ARG:HG2	1.88	0.73
2:B:286:TRP:NE1	2:B:290:LEU:HD12	2.05	0.72
1:A:300:ASN:CG	1:A:303:ILE:HD12	2.09	0.72
2:B:305:CYS:HB2	7:B:511:HOH:O	1.91	0.70
2:B:62:TYR:O	2:B:70:LYS:HE3	1.91	0.70
1:A:163:MET:SD	1:A:192:LYS:HG3	2.31	0.70
2:B:71:HIS:O	2:B:75:SER:HB3	1.92	0.70
1:A:300:ASN:OD1	1:A:303:ILE:HD12	1.92	0.69
1:A:214:CYS:HG	6:A:386:HG:HG	1.34	0.69
2:B:72:ILE:HG23	2:B:290:LEU:CD2	2.24	0.68
2:B:143:ASN:O	2:B:147:GLN:HG3	1.94	0.68
1:A:116:THR:O	1:A:120:ARG:HG3	1.94	0.67
2:B:258:ASP:OD2	2:B:261:MET:HB2	1.95	0.67
2:B:72:ILE:HG12	2:B:290:LEU:CD2	2.25	0.66
1:A:155:SER:O	1:A:159:GLU:HG3	1.95	0.66
1:A:7:GLN:NE2	7:A:545:HOH:O	2.27	0.66
2:B:163:MET:CE	2:B:189:LEU:HB2	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:ASP:CB	2:B:9:LYS:HE2	2.13	0.65
2:B:146:ILE:HD13	2:B:286:TRP:CZ3	2.31	0.65
2:B:300:ASN:OD1	2:B:302:ASP:N	2.30	0.65
2:B:291:PHE:HB3	2:B:301:LYS:HE3	1.77	0.65
2:B:306:GLN:NE2	2:B:306:GLN:HA	2.12	0.65
2:B:222:GLU:HA	2:B:225:GLU:OE2	1.97	0.64
2:B:19:PHE:CE1	2:B:98:LEU:HD22	2.32	0.63
2:B:31:GLN:NE2	7:B:417:HOH:O	2.30	0.63
1:A:205:ALA:O	1:A:315:ARG:HD2	1.99	0.63
2:B:149:ARG:O	2:B:282:GLN:NE2	2.30	0.63
1:A:286:TRP:NE1	1:A:290:LEU:HD22	2.14	0.63
2:B:267:GLU:HB3	7:B:533:HOH:O	1.99	0.63
2:B:284:LYS:HD3	7:B:511:HOH:O	1.98	0.62
2:B:323:LEU:HD23	2:B:323:LEU:N	2.12	0.62
2:B:315:ARG:HH11	2:B:315:ARG:HG2	1.64	0.62
1:A:286:TRP:HE1	1:A:290:LEU:HD22	1.65	0.62
2:B:127:ARG:NH1	7:B:510:HOH:O	2.29	0.62
1:A:217:ALA:HB2	1:A:299:LEU:HD13	1.80	0.62
2:B:217:ALA:HB2	2:B:299:LEU:HD22	1.81	0.62
2:B:149:ARG:NH1	2:B:286:TRP:NE1	2.48	0.62
2:B:315:ARG:HH11	2:B:315:ARG:CG	2.13	0.61
2:B:291:PHE:HB2	2:B:301:LYS:HG2	1.83	0.61
1:A:302:ASP:O	1:A:306:GLN:HG3	2.01	0.61
2:B:59:ARG:HB3	7:B:552:HOH:O	1.98	0.61
2:B:163:MET:HE2	2:B:189:LEU:CB	2.29	0.61
1:A:139:ASP:O	1:A:143:ASN:HB2	2.00	0.60
2:B:328:ARG:HB2	7:B:550:HOH:O	2.00	0.60
2:B:205:ALA:HB1	2:B:315:ARG:HG3	1.82	0.60
2:B:299:LEU:HD11	2:B:304:LEU:HD13	1.84	0.60
2:B:144:GLU:HB3	2:B:148:LYS:NZ	2.16	0.60
2:B:195:LEU:HD22	2:B:272[A]:CYS:SG	2.42	0.59
2:B:316:MET:HB3	2:B:321:LEU:HB2	1.85	0.59
2:B:149:ARG:NE	2:B:283:GLU:OE1	2.35	0.59
2:B:275:LEU:HD13	2:B:275:LEU:N	2.17	0.59
1:A:122:PHE:O	1:A:126:ILE:HG13	2.03	0.58
2:B:286:TRP:CE2	2:B:290:LEU:HD12	2.38	0.58
2:B:275:LEU:HD12	7:B:531:HOH:O	2.04	0.57
2:B:313:ASN:HD21	2:B:325:PHE:H	1.50	0.57
1:A:275:LEU:HD12	1:A:275:LEU:C	2.25	0.57
2:B:139:ASP:O	2:B:143:ASN:HB2	2.05	0.57
2:B:155:SER:O	2:B:159:GLU:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:LYS:NZ	1:A:258:ASP:OD2	2.29	0.56
2:B:52:GLU:HG2	2:B:52:GLU:O	2.06	0.56
1:A:284:LYS:NZ	1:A:309:GLU:HG3	2.21	0.56
2:B:205:ALA:HB2	2:B:242:LEU:HD21	1.87	0.56
1:A:312:THR:O	1:A:316:MET:HG3	2.06	0.55
2:B:72:ILE:CG1	2:B:290:LEU:HD23	2.33	0.55
2:B:145:GLN:HG3	2:B:289:TYR:CD2	2.42	0.55
1:A:49:ARG:HG3	7:A:524:HOH:O	2.06	0.54
2:B:116:THR:O	2:B:120:ARG:HG2	2.07	0.54
1:A:332:ILE:O	1:A:335:ILE:HG12	2.08	0.54
1:A:287:ALA:HB1	1:A:301:LYS:HD3	1.89	0.54
2:B:65:LEU:HD22	2:B:69:GLU:OE1	2.08	0.53
1:A:145:GLN:HG2	1:A:289:TYR:CD1	2.43	0.53
1:A:186:LEU:O	1:A:190:LYS:HG3	2.09	0.53
1:A:93:VAL:HG13	7:A:510:HOH:O	2.08	0.53
1:A:221:ARG:NH1	1:A:296:MET:HE2	2.24	0.53
1:A:253:ARG:NH1	1:A:266:GLU:OE1	2.42	0.52
2:B:278:GLN:NE2	2:B:278:GLN:HA	2.24	0.52
2:B:299:LEU:HG	2:B:304:LEU:HD22	1.92	0.51
2:B:117:ILE:HA	2:B:120:ARG:HG3	1.91	0.51
1:A:145:GLN:HG3	1:A:289:TYR:CB	2.40	0.51
2:B:149:ARG:NH1	2:B:286:TRP:CE2	2.79	0.50
1:A:31:GLN:NE2	7:A:401:HOH:O	2.29	0.50
2:B:222:GLU:O	2:B:222:GLU:OE2	2.29	0.50
2:B:219:ALA:HB2	2:B:228:ALA:HB2	1.94	0.50
1:A:222:GLU:C	1:A:223:LEU:HD23	2.33	0.50
2:B:190:LYS:HB2	2:B:264:ILE:HD13	1.94	0.49
1:A:287:ALA:HB3	1:A:301:LYS:CD	2.36	0.49
1:A:96:LEU:HB2	1:A:97:PRO:HD3	1.95	0.49
1:A:222:GLU:HG3	1:A:225:GLU:OE2	2.12	0.49
1:A:144:GLU:O	1:A:148:LYS:N	2.41	0.49
1:A:335:ILE:HG13	1:A:336:ASN:N	2.28	0.49
1:A:287:ALA:HB2	1:A:304:LEU:HD12	1.95	0.49
2:B:255:GLY:HA2	2:B:262:ALA:HB2	1.94	0.49
1:A:301:LYS:HD2	1:A:305[A]:CYS:HG	1.77	0.48
1:A:59:ARG:NH2	1:A:60:ILE:HD13	2.28	0.48
1:A:284:LYS:HZ1	1:A:309:GLU:CG	2.26	0.48
1:A:31:GLN:HG3	1:A:34:ASP:HA	1.94	0.48
2:B:291:PHE:O	2:B:301:LYS:HE3	2.13	0.48
2:B:310:TYR:CZ	2:B:330:ASN:HB2	2.48	0.48
1:A:268:CYS:O	1:A:271:GLU:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:LEU:HD21	2:B:268:CYS:SG	2.53	0.48
1:A:172:GLU:HG3	1:A:184:VAL:O	2.13	0.48
1:A:143:ASN:HB3	1:A:146:ILE:HG13	1.95	0.48
1:A:221:ARG:HH12	1:A:296:MET:HB3	1.78	0.48
1:A:203:LEU:HD23	1:A:203:LEU:C	2.34	0.48
2:B:300:ASN:HB2	7:B:464:HOH:O	2.14	0.47
2:B:242:LEU:HD13	2:B:242:LEU:C	2.35	0.47
1:A:327:THR:HG22	1:A:328:ARG:N	2.30	0.47
1:A:68:HIS:O	1:A:72:ILE:HG13	2.15	0.47
2:B:82:LEU:HD22	2:B:146:ILE:CG2	2.45	0.46
1:A:328:ARG:N	7:A:462:HOH:O	2.46	0.46
2:B:47:PHE:HE2	2:B:49:ARG:CZ	2.29	0.46
1:A:255:GLY:HA2	1:A:262:ALA:HB2	1.97	0.46
2:B:38:LYS:N	2:B:38:LYS:HD2	2.29	0.46
1:A:143:ASN:O	1:A:146:ILE:HB	2.15	0.46
1:A:317:GLN:OE1	1:A:323:LEU:HD11	2.15	0.46
2:B:273:TYR:CE1	2:B:277:VAL:HG21	2.51	0.46
1:A:300:ASN:ND2	1:A:303:ILE:HD12	2.30	0.46
2:B:124:HIS:O	2:B:128:ASN:ND2	2.29	0.46
2:B:37:GLU:OE2	2:B:41:GLU:OE2	2.33	0.46
1:A:63:GLN:HG2	1:A:64:ALA:N	2.31	0.46
1:A:163:MET:HE2	1:A:189:LEU:HA	1.98	0.46
1:A:74:ILE:CG2	1:A:75:SER:N	2.79	0.46
2:B:144:GLU:HB3	2:B:148:LYS:HZ2	1.79	0.45
1:A:96:LEU:N	1:A:97:PRO:CD	2.80	0.45
2:B:146:ILE:O	2:B:146:ILE:HG22	2.16	0.45
2:B:221:ARG:O	2:B:222:GLU:HB3	2.16	0.45
2:B:27:ARG:NH2	2:B:29:ASP:OD2	2.30	0.45
1:A:221:ARG:NH1	1:A:296:MET:CE	2.80	0.45
1:A:145:GLN:CG	1:A:289:TYR:CG	3.00	0.45
1:A:31:GLN:HG3	1:A:34:ASP:HB3	1.99	0.45
2:B:152:GLY:N	7:B:428:HOH:O	2.36	0.45
1:A:59:ARG:NH2	1:A:60:ILE:CD1	2.80	0.44
1:A:65:LEU:O	1:A:70:LYS:NZ	2.50	0.44
2:B:51:GLU:HB2	7:B:502:HOH:O	2.17	0.44
1:A:31:GLN:HG3	1:A:34:ASP:CB	2.46	0.44
1:A:130:VAL:CG1	1:A:131:ASN:N	2.80	0.44
2:B:92:ASN:O	2:B:96:LEU:HB2	2.17	0.44
2:B:79:TYR:CZ	2:B:149:ARG:HG2	2.53	0.44
1:A:307:TYR:CD1	1:A:332:ILE:HD11	2.52	0.44
1:A:307:TYR:CE1	1:A:332:ILE:HD11	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:LYS:HZ1	1:A:309:GLU:HG3	1.82	0.44
1:A:247:HIS:ND1	7:A:421:HOH:O	2.29	0.44
2:B:72:ILE:CG2	2:B:290:LEU:HD21	2.35	0.44
2:B:311:ILE:HB	2:B:315:ARG:NH1	2.33	0.44
2:B:63:GLN:HE21	2:B:63:GLN:CA	2.30	0.44
2:B:131:ASN:HD22	2:B:131:ASN:HA	1.56	0.43
1:A:219:ALA:HB1	1:A:338:TRP:CH2	2.52	0.43
1:A:1:ALA:HB3	7:A:417:HOH:O	2.17	0.43
2:B:275:LEU:HA	2:B:275:LEU:HD12	1.63	0.43
1:A:35:ILE:HG23	1:A:36:PHE:N	2.33	0.43
2:B:142:THR:O	2:B:147:GLN:OE1	2.37	0.43
2:B:145:GLN:HG3	2:B:289:TYR:CE2	2.53	0.43
1:A:221:ARG:HH12	1:A:296:MET:HE2	1.83	0.42
2:B:78:LYS:NZ	7:B:452:HOH:O	2.49	0.42
2:B:146:ILE:CD1	2:B:286:TRP:CH2	3.02	0.42
2:B:313:ASN:HD22	2:B:324:PRO:HD2	1.83	0.42
2:B:206:ILE:HB	2:B:311:ILE:CD1	2.49	0.42
1:A:223:LEU:N	1:A:223:LEU:HD23	2.34	0.42
1:A:65:LEU:HA	1:A:66:PRO:HD2	1.87	0.42
2:B:286:TRP:NE1	2:B:290:LEU:CD1	2.80	0.42
1:A:74:ILE:HG23	1:A:78:LYS:HE3	2.02	0.42
2:B:96:LEU:N	2:B:97:PRO:CD	2.83	0.42
2:B:127:ARG:CG	2:B:127:ARG:HH11	2.31	0.42
2:B:252:LEU:HD23	2:B:252:LEU:HA	1.89	0.42
1:A:145:GLN:HA	1:A:145:GLN:NE2	2.35	0.42
1:A:296:MET:HB3	1:A:296:MET:HE2	1.91	0.42
2:B:79:TYR:HE1	2:B:150:ALA:HB2	1.85	0.42
1:A:145:GLN:HG3	1:A:289:TYR:HB2	2.02	0.42
2:B:69:GLU:CD	2:B:221:ARG:HH22	2.21	0.41
2:B:69:GLU:HG2	2:B:296:MET:HG2	2.02	0.41
1:A:74:ILE:HD12	1:A:74:ILE:HA	1.77	0.41
2:B:311:ILE:HG13	2:B:312:THR:N	2.35	0.41
1:A:317:GLN:HG2	7:A:492:HOH:O	2.21	0.41
1:A:66:PRO:O	1:A:69:GLU:N	2.51	0.41
1:A:195:LEU:O	1:A:199:SER:OG	2.30	0.41
1:A:92:ASN:O	1:A:96:LEU:HB2	2.20	0.41
2:B:255:GLY:CA	2:B:262:ALA:HB2	2.50	0.41
2:B:79:TYR:HD1	2:B:146:ILE:HG23	1.86	0.41
2:B:221:ARG:HH11	2:B:221:ARG:HD2	1.71	0.41
2:B:105:GLU:O	2:B:109:GLU:HG3	2.21	0.41
2:B:301:LYS:HD2	7:B:464:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:120:ARG:HD2	2:B:120:ARG:HH11	1.65	0.41
1:A:96:LEU:HA	1:A:96:LEU:HD23	1.81	0.41
2:B:63:GLN:NE2	2:B:63:GLN:HA	2.36	0.40
2:B:260:GLU:O	2:B:264:ILE:HD12	2.22	0.40
1:A:149:ARG:NH2	1:A:283:GLU:OE1	2.45	0.40
2:B:149:ARG:NH2	2:B:283:GLU:OE1	2.53	0.40
1:A:301:LYS:HG2	1:A:301:LYS:HZ2	1.52	0.40
2:B:205:ALA:HB1	2:B:315:ARG:CG	2.48	0.40
2:B:197:LEU:HA	2:B:197:LEU:HD23	1.93	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	337/375 (90%)	327 (97%)	10 (3%)	0	100	100
2	B	339/375 (90%)	328 (97%)	11 (3%)	0	100	100
All	All	676/750 (90%)	655 (97%)	21 (3%)	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	305/338 (90%)	277 (91%)	28 (9%)	11 5
2	B	307/339 (91%)	274 (89%)	33 (11%)	8 2
All	All	612/677 (90%)	551 (90%)	61 (10%)	9 3

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	31	GLN
1	A	38	LYS
1	A	49	ARG
1	A	74	ILE
1	A	100	SER
1	A	130	VAL
1	A	145	GLN
1	A	148	LYS
1	A	155	SER
1	A	186	LEU
1	A	187	ARG
1	A	199	SER
1	A	222	GLU
1	A	253	ARG
1	A	263	GLU
1	A	271	GLU
1	A	292	ARG
1	A	295	SER
1	A	299	LEU
1	A	301	LYS
1	A	302	ASP
1	A	311	ILE
1	A	315	ARG
1	A	329	SER
1	A	337	THR
1	A	339	LEU
1	A	340	VAL
2	B	31	GLN
2	B	38	LYS
2	B	42	LYS
2	B	52	GLU
2	B	100	SER
2	B	120	ARG
2	B	127	ARG

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Mol	Chain	Res	Type
2	B	131	ASN
2	B	134	SER
2	B	148	LYS
2	B	155	SER
2	B	170	LEU
2	B	186	LEU
2	B	199	SER
2	B	206	ILE
2	B	208	PHE
2	B	222	GLU
2	B	232	ARG
2	B	254	SER
2	B	268	CYS
2	B	275	LEU
2	B	276	PHE
2	B	284	LYS
2	B	286	TRP
2	B	290	LEU
2	B	292	ARG
2	B	296	MET
2	B	297	ILE
2	B	304	LEU
2	B	313	ASN
2	B	315	ARG
2	B	317	GLN
2	B	329	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	71	HIS
1	A	246	GLN
1	A	278	GLN
1	A	281	GLN
1	A	306	GLN
2	B	63	GLN
2	B	80	GLN
2	B	131	ASN
2	B	201	ASN
2	B	227	ASN
2	B	247	HIS

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Mol	Chain	Res	Type
2	B	270	GLN
2	B	278	GLN
2	B	313	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	MTY	A	208	1	11,12,13	1.27	3 (27%)	12,15,17	1.23	1 (8%)

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
1	MTY	A	208	1	-	0/4/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	208	MTY	CB-CA	-2.14	1.49	1.53
1	A	208	MTY	CZ-CE2	-2.08	1.34	1.38
1	A	208	MTY	CD2-CG	-2.02	1.35	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	208	MTY	O-C-CA	-2.05	120.15	125.49

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 23 ligands modelled in this entry, 3 are modelled with single atom and 20 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.

No monomer is involved in short contacts.

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	339/375 (90%)	-0.16	3 (0%) 85 89	13, 26, 54, 75	0
2	B	341/375 (90%)	-0.02	13 (3%) 44 50	12, 28, 61, 86	1 (0%)
All	All	680/750 (90%)	-0.09	16 (2%) 62 68	12, 27, 57, 86	1 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	292	ARG	4.7
2	B	289	TYR	4.6
1	A	340	VAL	4.5
2	B	294	GLY	4.0
2	B	222	GLU	3.2
1	A	292	ARG	2.9
2	B	255	GLY	2.9
2	B	267	GLU	2.9
2	B	296	MET	2.8
2	B	148	LYS	2.8
2	B	293	ASP	2.7
2	B	295	SER	2.6
2	B	297	ILE	2.5
1	A	337	THR	2.5
2	B	259	PRO	2.2
2	B	151	GLU	2.0

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

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
1	MTY	A	208	12/13	0.90	0.17	-	22,29,58,100	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. 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
5	OH	B	378	1/1	0.97	0.16	1.44	17,17,17,17	0
5	OH	A	379	1/1	0.89	0.14	1.13	28,28,28,28	0
6	HG	B	382	1/1	0.94	0.14	-0.26	41,41,41,41	1
6	HG	A	386	1/1	0.96	0.12	-0.55	62,62,62,62	1
6	HG	B	385	1/1	0.96	0.11	-0.61	31,31,31,31	1
6	HG	B	381	1/1	0.99	0.08	-0.92	22,22,22,22	1
3	FE2	B	377	1/1	0.99	0.07	-1.16	24,24,24,24	0
5	OH	A	378	1/1	0.98	0.07	-1.22	15,15,15,15	0
6	HG	B	383	1/1	0.99	0.10	-1.24	32,32,32,32	1
6	HG	A	384	1/1	0.97	0.07	-1.32	34,34,34,34	1
6	HG	B	379	1/1	1.00	0.06	-1.54	20,20,20,20	1
4	FE	A	377	1/1	0.99	0.06	-2.06	20,20,20,20	0
6	HG	A	380	1/1	1.00	0.05	-2.31	23,23,23,23	1
3	FE2	B	376	1/1	0.99	0.05	-2.38	28,28,28,28	0
6	HG	A	381	1/1	0.98	0.06	-2.56	24,24,24,24	1
6	HG	B	386	1/1	0.99	0.08	-3.19	24,24,24,24	1
6	HG	A	382	1/1	1.00	0.05	-3.23	27,27,27,27	1
6	HG	A	385	1/1	0.99	0.03	-3.53	31,31,31,31	1
6	HG	B	387	1/1	0.99	0.07	-3.61	25,25,25,25	1
6	HG	B	380	1/1	1.00	0.04	-5.54	24,24,24,24	1
6	HG	B	384	1/1	0.99	0.07	-	24,24,24,24	1
6	HG	A	383	1/1	0.99	0.06	-	23,23,23,23	1
3	FE2	A	376	1/1	0.99	0.03	-	27,27,27,27	0

6.5 Other polymers (i)

There are no such residues in this entry.