



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:11 AM UTC

PDB ID : 1BIQ / pdb_00001biq
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

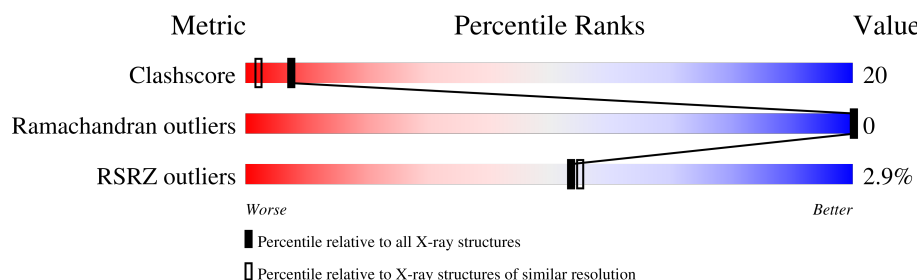
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	375	5% 58% 30% • 9%
2	B	375	5% 58% 29% • 9%

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 mutation	UNP P69924
A	208	MTY	PHE	modified residue	UNP P69924
A	238	ALA	GLU	engineered mutation	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 mutation	UNP P69924
B	238	ALA	GLU	engineered mutation	UNP P69924

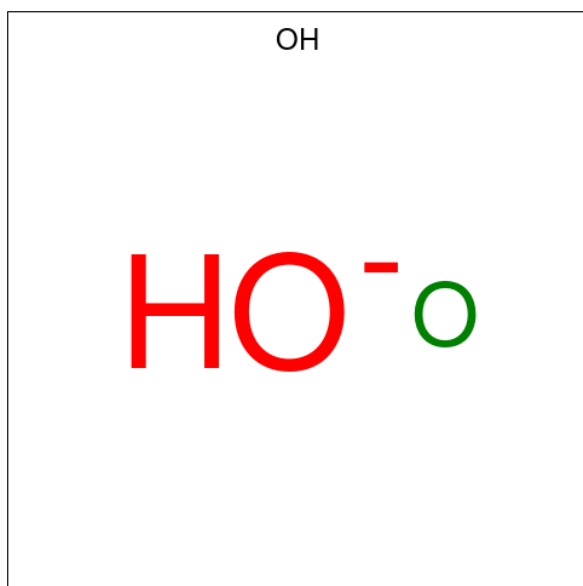
- Molecule 3 is FE (II) ION (CCD ID: FE2) (formula: Fe).

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

- Molecule 4 is FE (III) ION (CCD ID: FE) (formula: Fe).

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

- Molecule 5 is HYDROXIDE ION (CCD ID: OH) (formula: HO).



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

- Molecule 6 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

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

- Molecule 7 is water.

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

Continued on next page...

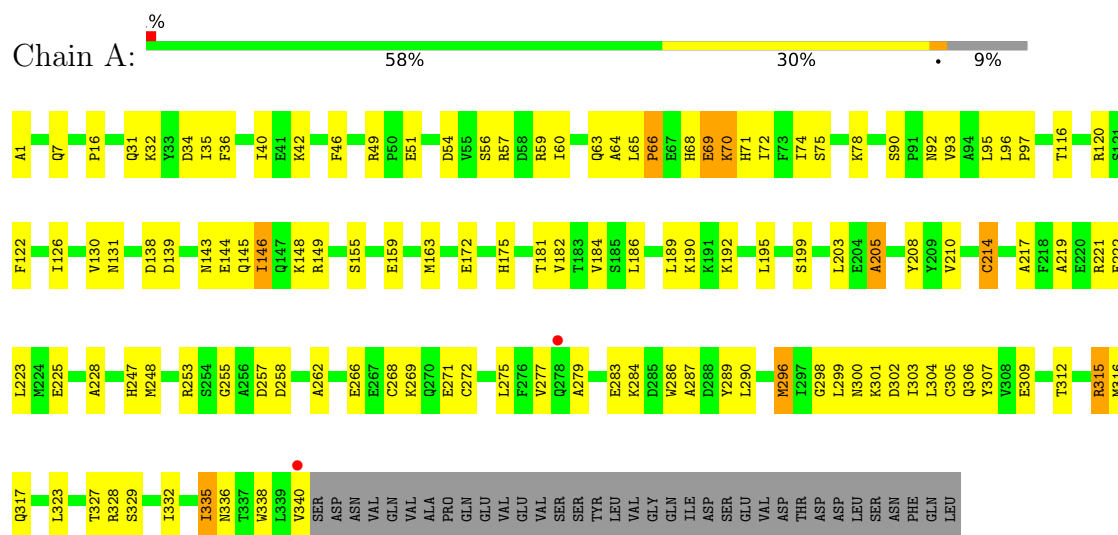
Continued from previous page...

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

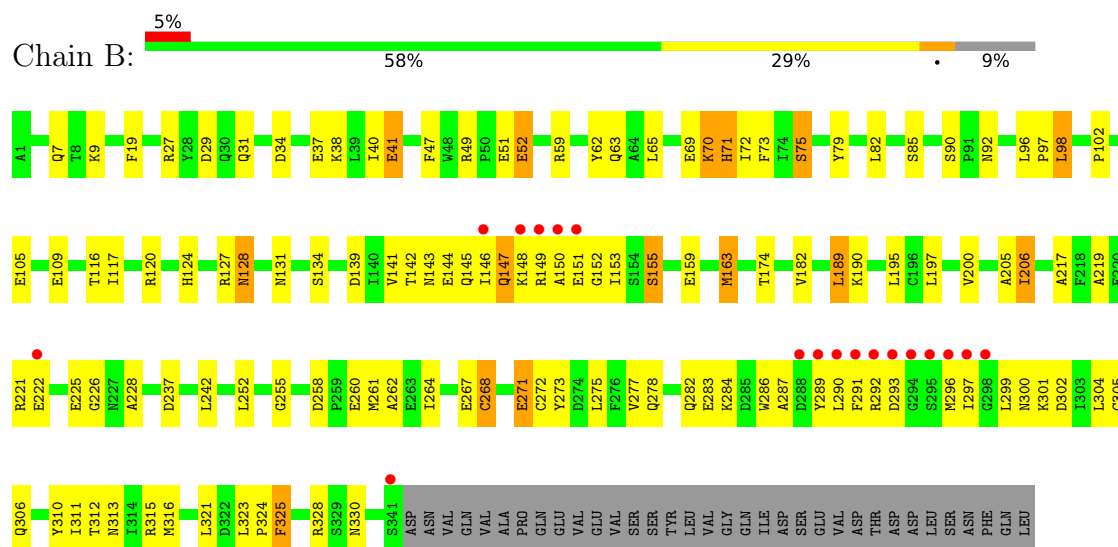
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE



• Molecule 2: PROTEIN R2 OF RIBONUCLEOTIDE REDUCTASE



4 Data and refinement statistics

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 25.00 – 2.06	Depositor EDS
% Data completeness (in resolution range)	98.0 (25.00-2.05) 98.1 (25.00-2.06)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.32 (at 2.06Å)	Xtriage
Refinement program	TNT 5E	Depositor
R, R_{free}	0.191 , 0.260 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.456	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 107.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MTY, HG, FE, FE2, 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	1.00	2/2834 (0.1%)	1.54	35/3843 (0.9%)
2	B	0.96	2/2853 (0.1%)	1.57	35/3870 (0.9%)
All	All	0.98	4/5687 (0.1%)	1.56	70/7713 (0.9%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	163	MET	SD-CE	-6.62	1.62	1.79
2	B	182	VAL	CA-CB	-6.40	1.46	1.54
1	A	269	LYS	CE-NZ	5.70	1.66	1.49
1	A	181	THR	CA-CB	5.30	1.60	1.53

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	206	ILE	N-CA-C	14.23	122.86	111.62
2	B	90	SER	N-CA-C	9.26	125.29	113.25
1	A	32	LYS	N-CA-C	-9.19	100.39	111.69
2	B	292	ARG	N-CA-C	8.98	123.55	112.23
2	B	226	GLY	N-CA-C	-8.78	101.87	112.49
2	B	293	ASP	N-CA-C	-8.24	102.94	113.16
2	B	271	GLU	N-CA-C	-8.18	102.30	111.14
1	A	329	SER	N-CA-C	-8.11	99.60	110.55
2	B	147	GLN	N-CA-C	-7.75	102.80	111.71
1	A	214	CYS	CB-CA-C	7.65	122.89	110.88
1	A	70	LYS	N-CA-C	-7.33	103.22	111.14
1	A	40	ILE	CB-CA-C	-7.21	102.74	111.97
2	B	287	ALA	N-CA-C	-6.94	103.72	111.71
1	A	90	SER	N-CA-C	6.92	122.24	113.25
2	B	189	LEU	N-CA-C	-6.68	103.48	111.69

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	VAL	CG1-CB-CG2	6.42	124.93	110.80
2	B	70	LYS	N-CA-C	-6.36	104.27	111.07
1	A	205	ALA	N-CA-C	6.31	121.05	113.23
2	B	153	ILE	N-CA-C	-6.28	105.03	110.74
1	A	95	LEU	N-CA-C	6.14	120.99	112.45
2	B	325	PHE	N-CA-C	6.04	118.39	109.69
2	B	41	GLU	N-CA-C	5.93	117.75	111.28
1	A	248	MET	N-CA-C	-5.79	104.87	111.07
1	A	257	ASP	CA-C-O	5.73	121.26	117.94
2	B	134	SER	CA-C-N	-5.68	113.40	120.56
2	B	134	SER	C-N-CA	-5.68	113.40	120.56
2	B	71	HIS	N-CA-C	5.64	117.11	111.07
2	B	128	ASN	N-CA-CB	-5.63	102.15	110.49
1	A	214	CYS	CA-C-N	5.62	127.74	120.44
1	A	214	CYS	C-N-CA	5.62	127.74	120.44
1	A	277	VAL	CA-C-N	-5.61	112.32	120.29
1	A	277	VAL	C-N-CA	-5.61	112.32	120.29
2	B	200	VAL	N-CA-C	-5.59	103.94	111.44
1	A	315	ARG	NE-CZ-NH1	5.59	127.09	121.50
1	A	210	VAL	CB-CA-C	-5.58	104.52	112.22
1	A	16	PRO	N-CA-C	-5.57	103.42	111.22
1	A	298	GLY	CA-C-N	-5.52	113.13	121.86
1	A	298	GLY	C-N-CA	-5.52	113.13	121.86
1	A	210	VAL	N-CA-C	-5.50	105.16	110.72
1	A	66	PRO	N-CA-C	-5.48	102.58	111.19
1	A	272	CYS	N-CA-C	-5.47	105.00	110.97
1	A	182	VAL	N-CA-C	-5.43	100.30	108.12
2	B	98	LEU	N-CA-C	5.42	120.56	113.30
1	A	71	HIS	CB-CA-C	-5.39	101.52	110.68
2	B	151	GLU	N-CA-C	5.37	117.81	110.24
1	A	296	MET	N-CA-C	-5.36	100.92	109.23
1	A	146	ILE	CB-CA-C	-5.35	105.12	111.97
2	B	73	PHE	N-CA-C	5.34	116.79	110.97
1	A	279	ALA	N-CA-C	-5.33	105.37	111.07
1	A	228	ALA	N-CA-C	5.28	117.78	111.71
2	B	174	THR	N-CA-C	-5.27	99.81	108.41
2	B	102	PRO	N-CA-C	5.26	120.84	113.53
2	B	268	CYS	N-CA-C	5.26	119.67	112.88
1	A	90	SER	CA-C-O	5.26	123.47	118.63
2	B	52	GLU	CA-C-N	-5.25	116.67	123.19
2	B	52	GLU	C-N-CA	-5.25	116.67	123.19
1	A	69	GLU	N-CA-C	-5.25	105.62	112.23

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	75	SER	N-CA-C	5.23	116.98	111.28
1	A	57	ARG	N-CA-C	5.22	116.97	111.28
2	B	40	ILE	CB-CA-C	-5.21	105.30	111.97
2	B	237	ASP	N-CA-C	-5.21	105.68	111.36
1	A	335	ILE	CB-CA-C	-5.20	103.68	112.16
2	B	141	VAL	N-CA-C	5.20	115.42	110.53
2	B	206	ILE	N-CA-CB	5.19	115.90	111.64
2	B	155	SER	N-CA-C	5.19	117.84	111.82
2	B	73	PHE	CB-CA-C	-5.18	102.98	110.96
1	A	175	HIS	N-CA-CB	-5.14	103.25	111.43
1	A	269	LYS	N-CA-C	5.13	121.73	110.80
2	B	85	SER	N-CA-C	-5.07	105.64	111.07
2	B	7	GLN	N-CA-C	5.06	119.32	113.20

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	124	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	0	0
7	A	160	0	0	9	0
7	B	165	0	0	13	0
All	All	5921	0	5461	218	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 (218) 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
1:A:221:ARG:NH1	1:A:296:MET:HE3	1.87	0.90
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
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: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:300:ASN:CG	1:A:303:ILE:HD12	2.09	0.76
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: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
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
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
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:222:GLU:HA	2:B:225:GLU:OE2	1.97	0.64
2:B:306:GLN:HA	2:B:306:GLN:NE2	2.12	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:163:MET:HE2	2:B:189:LEU:HB2	1.79	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:N	2:B:323:LEU:HD23	2.12	0.62
1:A:286:TRP:HE1	1:A:290:LEU:HD22	1.65	0.62
2:B:163:MET:HE3	7:B:442:HOH:O	2.00	0.62
2:B:315:ARG:HG2	2:B:315:ARG:HH11	1.64	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
1:A:275:LEU:HD12	1:A:275:LEU:C	2.25	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
1:A:302:ASP:O	1:A:306:GLN:HG3	2.01	0.61
2:B:291:PHE:HB2	2:B:301:LYS:HG2	1.83	0.61
2:B:59:ARG:HB3	7:B:552:HOH:O	1.98	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:144:GLU:HB3	2:B:148:LYS:NZ	2.16	0.60
2:B:299:LEU:HD11	2:B:304:LEU:HD13	1.84	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:N	2:B:275:LEU:HD13	2.17	0.59
1:A:122:PHE:O	1:A:126:ILE:HG13	2.03	0.58
1:A:221:ARG:CZ	1:A:296:MET:HE3	2.33	0.58
2:B:286:TRP:CE2	2:B:290:LEU:HD12	2.38	0.58
1:A:284:LYS:HZ1	1:A:309:GLU:CG	2.17	0.57
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
2:B:139:ASP:O	2:B:143:ASN:HB2	2.05	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:155:SER:O	2:B:159:GLU:HG3	2.05	0.56
2:B:52:GLU:HG2	2:B:52:GLU:O	2.06	0.56
2:B:217:ALA:CB	2:B:296:MET:HE1	2.35	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:190:LYS:NZ	1:A:258:ASP:OD2	2.29	0.55
1:A:49:ARG:HG3	7:A:524:HOH:O	2.06	0.54
1:A:222:GLU:C	1:A:223:LEU:HD23	2.33	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
1:A:284:LYS:HZ1	1:A:309:GLU:HG3	1.73	0.54
2:B:65:LEU:HD22	2:B:69:GLU:OE1	2.08	0.53
2:B:217:ALA:HB3	2:B:296:MET:HE1	1.91	0.53
1:A:145:GLN:HG2	1:A:289:TYR:CD1	2.43	0.53
1:A:93:VAL:HG13	7:A:510:HOH:O	2.08	0.53
1:A:186:LEU:O	1:A:190:LYS:HG3	2.09	0.53
1:A:203:LEU:C	1:A:203:LEU:HD23	2.34	0.52
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
1:A:221:ARG:HH12	1:A:296:MET:HE3	1.73	0.52
2:B:163:MET:HE1	2:B:189:LEU:CA	2.40	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
2:B:242:LEU:C	2:B:242:LEU:HD13	2.35	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
1:A:301:LYS:HD2	1:A:305[A]:CYS:HG	1.74	0.50
2:B:219:ALA:HB2	2:B:228:ALA:HB2	1.94	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:144:GLU:O	1:A:148:LYS:N	2.41	0.49
1:A:222:GLU:HG3	1:A:225:GLU:OE2	2.12	0.49
1:A:335:ILE:HG13	1:A:336:ASN:N	2.28	0.49
1:A:54:ASP:OD1	1:A:56:SER:OG	2.29	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:59:ARG:NH2	1:A:60:ILE:HD13	2.28	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
1:A:268:CYS:O	1:A:271:GLU:HB2	2.14	0.48
2:B:310:TYR:CZ	2:B:330:ASN:HB2	2.48	0.48
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:338:TRP:H	1:A:338:TRP:CD1	2.31	0.47
2:B:163:MET:HE1	2:B:189:LEU:N	2.29	0.47
2:B:300:ASN:HB2	7:B:464:HOH:O	2.14	0.47
1:A:68:HIS:O	1:A:72:ILE:HG13	2.15	0.47
1:A:327:THR:HG22	1:A:328:ARG:N	2.30	0.47
1:A:328:ARG:N	7:A:462:HOH:O	2.46	0.46
2:B:82:LEU:HD22	2:B:146:ILE:CG2	2.45	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
2:B:47:PHE:HE2	2:B:49:ARG:CZ	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:37:GLU:OE2	2:B:41:GLU:OE2	2.33	0.46
2:B:69:GLU:CD	2:B:221:ARG:HH22	2.21	0.46
2:B:124:HIS:O	2:B:128:ASN:ND2	2.29	0.46
2:B:144:GLU:HB3	2:B:148:LYS:HZ1	1.81	0.46
1:A:74:ILE:CG2	1:A:75:SER:N	2.79	0.46
1:A:214:CYS:HG	6:A:386:HG:HG	1.53	0.46
1:A:96:LEU:N	1:A:97:PRO:CD	2.80	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
2:B:163:MET:HE2	2:B:189:LEU:CB	2.46	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
1:A:63:GLN:HG2	1:A:64:ALA:N	2.31	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:A:247:HIS:ND1	7:A:421:HOH:O	2.29	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
2:B:79:TYR:CZ	2:B:149:ARG:HG2	2.53	0.44
2:B:72:ILE:CG2	2:B:290:LEU:HD21	2.35	0.44
2:B:63:GLN:CA	2:B:63:GLN:HE21	2.30	0.44
2:B:311:ILE:HB	2:B:315:ARG:NH1	2.33	0.44
2:B:131:ASN:HD22	2:B:131:ASN:HA	1.56	0.43
2:B:146:ILE:O	2:B:146:ILE:HG22	2.16	0.43
2:B:296:MET:HE3	2:B:296:MET:HB3	1.77	0.43
1:A:1:ALA:HB3	7:A:417:HOH:O	2.17	0.43
1:A:163:MET:HE2	1:A:189:LEU:HA	2.00	0.43
1:A:219:ALA:HB1	1:A:338:TRP:CH2	2.52	0.43
2:B:275:LEU:HD12	2:B:275:LEU:HA	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
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
1:A:65:LEU:HA	1:A:66:PRO:HD2	1.87	0.42
1:A:223:LEU:HD23	1:A:223:LEU:N	2.34	0.42
2:B:206:ILE:HB	2:B:311:ILE:CD1	2.49	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
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
1:A:74:ILE:HA	1:A:74:ILE:HD12	1.77	0.41
2:B:69:GLU:HG2	2:B:296:MET:HG2	2.02	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:105:GLU:O	2:B:109:GLU:HG3	2.21	0.41
2:B:221:ARG:HH11	2:B:221:ARG:HD2	1.71	0.41
1:A:96:LEU:HA	1:A:96:LEU:HD23	1.81	0.41
2:B:120:ARG:HH11	2:B:120:ARG:HD2	1.65	0.41
2:B:301:LYS:HD2	7:B:464:HOH:O	2.21	0.41
1:A:51:GLU:H	1:A:51:GLU:CD	2.29	0.40
2:B:63:GLN:NE2	2:B:63:GLN:HA	2.36	0.40
1:A:149:ARG:NH2	1:A:283:GLU:OE1	2.45	0.40
2:B:260:GLU:O	2:B:264:ILE:HD12	2.22	0.40
2:B:311:ILE:HG13	2:B:312:THR:N	2.35	0.40
2:B:149:ARG:NH2	2:B:283:GLU:OE1	2.53	0.40
2:B:197:LEU:HD23	2:B:197:LEU:HA	1.93	0.40
2:B:205:ALA:HB1	2:B:315:ARG:CG	2.48	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](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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	1.35	3 (27%)	10,15,17	1.18	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
1	MTY	A	208	1	-	4/5/6/8	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	CZ-CE2	-2.31	1.34	1.39
1	A	208	MTY	CB-CA	-2.16	1.49	1.53
1	A	208	MTY	CD2-CG	-2.14	1.35	1.39

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	208	MTY	C-CA-CB-CG
1	A	208	MTY	CA-CB-CG-CD2
1	A	208	MTY	N-CA-CB-CG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
1	A	208	MTY	CA-CB-CG-CD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 20 are monoatomic and 3 are modelled with single atom - 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, 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	339/375 (90%)	-0.06	2 (0%) 85 87	13, 26, 54, 75	0
2	B	341/375 (90%)	0.04	18 (5%) 32 32	12, 28, 61, 86	1 (0%)
All	All	680/750 (90%)	-0.01	20 (2%) 53 55	12, 27, 57, 86	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	294	GLY	4.1
2	B	296	MET	3.6
2	B	289	TYR	3.6
1	A	340	VAL	3.2
2	B	148	LYS	3.0
2	B	291	PHE	2.9
2	B	150	ALA	2.8
2	B	295	SER	2.8
2	B	222	GLU	2.8
2	B	297	ILE	2.7
2	B	290	LEU	2.6
1	A	278	GLN	2.5
2	B	146	ILE	2.4
2	B	292	ARG	2.3
2	B	149	ARG	2.2
2	B	151	GLU	2.2
2	B	293	ASP	2.2
2	B	288	ASP	2.1
2	B	298	GLY	2.0
2	B	341	SER	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	MTY	A	208	12/13	0.90	0.10	22,29,58,100	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	OH	A	379	1/1	0.89	0.13	28,28,28,28	0
6	HG	A	386	1/1	0.89	0.12	62,62,62,62	1
6	HG	B	382	1/1	0.95	0.10	41,41,41,41	1
5	OH	A	378	1/1	0.96	0.06	15,15,15,15	0
6	HG	A	384	1/1	0.96	0.06	34,34,34,34	1
6	HG	B	385	1/1	0.96	0.07	31,31,31,31	1
6	HG	A	381	1/1	0.97	0.04	24,24,24,24	1
6	HG	B	384	1/1	0.98	0.07	24,24,24,24	1
5	OH	B	378	1/1	0.98	0.04	17,17,17,17	0
6	HG	B	386	1/1	0.98	0.05	24,24,24,24	1
3	FE2	A	376	1/1	0.99	0.02	27,27,27,27	0
6	HG	A	385	1/1	0.99	0.03	31,31,31,31	1
3	FE2	B	376	1/1	0.99	0.02	28,28,28,28	0
6	HG	B	381	1/1	0.99	0.05	22,22,22,22	1
3	FE2	B	377	1/1	0.99	0.02	24,24,24,24	0
6	HG	B	383	1/1	0.99	0.07	32,32,32,32	1
4	FE	A	377	1/1	0.99	0.03	20,20,20,20	0
6	HG	A	382	1/1	0.99	0.03	27,27,27,27	1
6	HG	A	383	1/1	0.99	0.03	23,23,23,23	1
6	HG	B	387	1/1	0.99	0.04	25,25,25,25	1
6	HG	B	379	1/1	1.00	0.02	20,20,20,20	1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	HG	B	380	1/1	1.00	0.02	24,24,24,24	1
6	HG	A	380	1/1	1.00	0.02	23,23,23,23	1

6.5 Other polymers [i](#)

There are no such residues in this entry.