



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:08 PM GMT

PDB ID : 3T1A
Title : Crystal Structure of HIV-1 Reverse Transcriptase (K103N mutant) in Complex with Inhibitor M05
Authors : Yan, Y.; Reid, J.
Deposited on : 2011-07-21
Resolution : 2.40 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

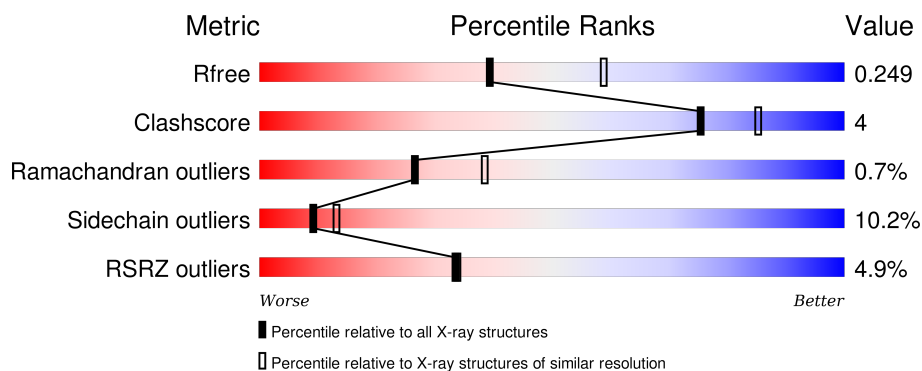
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	563	 3% 82% 15% ..
1	B	563	 5% 57% 12% • 29%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8113 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.

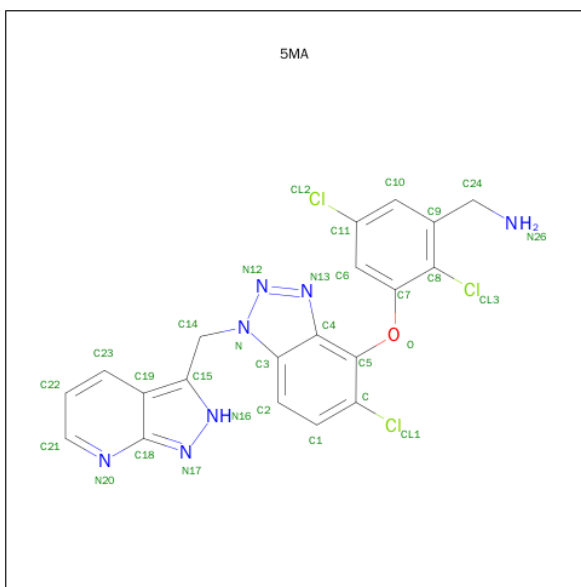
- Molecule 1 is a protein called Reverse Transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	0	0
			4503	2911	751	833	8			
1	B	400	Total	C	N	O	S	0	0	0
			3311	2155	548	602	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	EXPRESSION TAG	UNP P04585
A	-1	ASN	-	EXPRESSION TAG	UNP P04585
A	0	SER	-	EXPRESSION TAG	UNP P04585
A	103	ASN	LYS	ENGINEERED MUTATION	UNP P04585
B	-2	MET	-	EXPRESSION TAG	UNP P04585
B	-1	ASN	-	EXPRESSION TAG	UNP P04585
B	0	SER	-	EXPRESSION TAG	UNP P04585
B	103	ASN	LYS	ENGINEERED MUTATION	UNP P04585

- Molecule 2 is 1-(2,5-DICHLORO-3-{[5-CHLORO-1-(2H-PYRAZOLO[3,4-B]PYRIDIN-3-YLMETHYL)-1H-BENZOTRIAZOL-4-YL]OXY}PHENYL)METHANAMINE (three-letter code: 5MA) (formula: C₂₀H₁₄Cl₃N₇O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			31	20	3	7	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	159	Total	O	0	0
			159	159		
3	B	109	Total	O	0	0
			109	109		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.33Å 154.23Å 155.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.14 – 2.40 34.90 – 2.40	Depositor EDS
% Data completeness (in resolution range)	96.8 (35.14-2.40) 96.9 (34.90-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.39Å)	Xtriage
Refinement program	BUSTER 2.9.7	Depositor
R, R_{free}	0.205 , 0.250 0.205 , 0.249	Depositor DCC
R_{free} test set	2753 reflections (5.37%)	DCC
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 54067 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8113	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 5MA

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	0.50	0/4619	0.75	1/6279 (0.0%)
1	B	0.50	0/3404	0.72	0/4627
All	All	0.50	0/8023	0.74	1/10906 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	345	PRO	C-N-CA	7.46	140.34	121.70

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	4503	0	4540	31	0
1	B	3311	0	3326	31	0
2	A	31	0	14	2	0
3	A	159	0	0	4	0
3	B	109	0	0	1	0
All	All	8113	0	7880	61	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 (61) 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:330:GLN:HE22	1:A:340:GLN:HE22	1.28	0.77
1:B:317:VAL:HG13	1:B:347:LYS:HG2	1.74	0.67
1:B:337:TRP:HE1	1:B:367:GLN:HE21	1.45	0.64
1:B:373:GLN:O	1:B:377:THR:HG23	1.98	0.64
1:A:195:ILE:HD12	1:A:199:ARG:HE	1.63	0.63
1:A:103:ASN:H	2:A:561:5MA:HN16	1.47	0.63
1:A:244:ILE:HD13	1:A:267:ALA:HB2	1.83	0.60
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.84	0.60
1:B:214:LEU:HD23	1:B:214:LEU:H	1.68	0.59
1:B:394:GLN:HB3	1:B:397:THR:HG23	1.84	0.58
1:A:175:ASN:HD21	1:A:201:LYS:HZ2	1.52	0.57
1:B:275:LYS:H	1:B:306:ASN:HD21	1.50	0.57
1:A:59:PRO:HG2	1:A:76:ASP:HB3	1.89	0.53
1:A:175:ASN:HD21	1:A:201:LYS:NZ	2.06	0.53
1:A:406:TRP:CH2	1:B:418:ASN:HA	2.44	0.53
1:A:318:TYR:CZ	2:A:561:5MA:H14A	2.45	0.52
1:A:60:VAL:HG12	1:A:75:VAL:HG22	1.91	0.51
1:B:335:GLY:HA3	1:B:356:ARG:HG2	1.92	0.51
1:B:180:ILE:HG12	1:B:189:VAL:HG13	1.91	0.51
1:A:399:GLU:HA	1:A:402:TRP:CE3	2.46	0.50
1:A:193:LEU:HB3	1:A:197:GLN:HG3	1.94	0.49
1:B:317:VAL:CG1	1:B:347:LYS:HG2	2.43	0.49
1:A:7:THR:HG21	3:A:597:HOH:O	2.12	0.49
1:B:344:GLU:H	1:B:347:LYS:HD3	1.78	0.48
1:B:115:TYR:HB3	1:B:149:LEU:HB2	1.95	0.48
1:B:163:SER:HA	1:B:166:LYS:HE3	1.95	0.48
1:A:395:LYS:NZ	1:A:416:PHE:H	2.12	0.47
1:B:287:LYS:HE3	1:B:293:ILE:HG21	1.97	0.47
1:B:343:GLN:H	1:B:347:LYS:HE2	1.80	0.46
1:B:345:PRO:O	1:B:346:PHE:HB2	2.15	0.46
1:B:244:ILE:HB	1:B:310:LEU:HG	1.98	0.46
1:B:373:GLN:NE2	1:B:407:GLN:H	2.13	0.46
1:A:296:THR:HG21	3:A:566:HOH:O	2.16	0.45
1:B:373:GLN:HE22	1:B:407:GLN:H	1.65	0.45
1:B:57:ASN:OD1	1:B:143:ARG:NH1	2.50	0.44
1:B:253:THR:HG22	1:B:256:ASP:H	1.82	0.44
1:B:269:GLN:HB3	1:B:346:PHE:CE1	2.53	0.44
1:A:365:VAL:O	1:A:369:THR:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:TRP:HZ2	1:B:427:TYR:CZ	2.35	0.43
1:A:57:ASN:HD22	1:A:143:ARG:NH1	2.15	0.43
1:B:377:THR:HG22	1:B:410:TRP:HZ2	1.82	0.43
1:A:175:ASN:HB3	1:A:178:ILE:HD12	2.01	0.43
1:B:325:LEU:HD12	1:B:385:LYS:HG3	2.00	0.43
1:A:184:MET:HE2	3:A:601:HOH:O	2.18	0.43
1:B:282:LEU:HB3	1:B:293:ILE:HD11	2.00	0.43
1:A:246:LEU:O	1:A:307:ARG:NH2	2.52	0.43
1:A:5:ILE:HG22	1:A:212:TRP:CE3	2.54	0.42
1:B:111:VAL:HG21	1:B:187:LEU:HD22	2.00	0.42
1:A:181:TYR:HB2	1:A:188:TYR:HB2	2.01	0.42
1:A:57:ASN:HD22	1:A:143:ARG:HH12	1.66	0.42
1:A:288:ALA:O	1:A:291:GLU:HB3	2.20	0.42
1:B:401:TRP:HE3	1:B:404:GLU:HG3	1.85	0.41
1:B:58:THR:HG23	1:B:76:ASP:O	2.20	0.41
1:A:296:THR:HG23	1:A:299:ALA:H	1.85	0.41
1:A:63:ILE:HG13	1:A:72:ARG:HG3	2.03	0.41
1:B:79:GLU:HG3	1:B:83:ARG:HE	1.86	0.41
1:A:21:VAL:HG22	1:A:59:PRO:HD3	2.02	0.41
1:B:172:ARG:HD2	3:B:616:HOH:O	2.20	0.41
1:A:369:THR:HG22	1:A:411:ILE:HD11	2.03	0.40
1:A:255:ASN:HB2	1:A:289:LEU:HD23	2.02	0.40
1:A:395:LYS:HE3	3:A:574:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/563 (98%)	525 (96%)	20 (4%)	5 (1%)	21	30
1	B	392/563 (70%)	379 (97%)	11 (3%)	2 (0%)	34	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	942/1126 (84%)	904 (96%)	31 (3%)	7 (1%)	26	38

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	334	GLN
1	B	347	LYS
1	A	543	GLY
1	A	491	LEU
1	B	346	PHE
1	A	195	ILE
1	A	219	LYS

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	492/503 (98%)	443 (90%)	49 (10%)	9	14
1	B	364/503 (72%)	326 (90%)	38 (10%)	9	12
All	All	856/1006 (85%)	769 (90%)	87 (10%)	9	13

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	6	GLU
1	A	7	THR
1	A	21	VAL
1	A	22	LYS
1	A	26	LEU
1	A	64	LYS
1	A	69	THR
1	A	70	LYS
1	A	91	GLN
1	A	94	ILE

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Mol	Chain	Res	Type
1	A	118	VAL
1	A	126	LYS
1	A	138	GLU
1	A	161	GLN
1	A	173	LYS
1	A	194	GLU
1	A	195	ILE
1	A	205	LEU
1	A	211	ARG
1	A	219	LYS
1	A	220	LYS
1	A	248	GLU
1	A	260	LEU
1	A	276	VAL
1	A	279	LEU
1	A	287	LYS
1	A	289	LEU
1	A	297	GLU
1	A	303	LEU
1	A	311	LYS
1	A	357	MET
1	A	368	LEU
1	A	373	GLN
1	A	374	LYS
1	A	402	TRP
1	A	428	GLN
1	A	449	GLU
1	A	459	THR
1	A	463	ARG
1	A	466	VAL
1	A	479	LEU
1	A	491	LEU
1	A	493	VAL
1	A	496	VAL
1	A	503	LEU
1	A	528	LYS
1	A	547	GLN
1	A	548	VAL
1	B	6	GLU
1	B	11	LYS
1	B	17	ASP
1	B	26	LEU

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Mol	Chain	Res	Type
1	B	42	GLU
1	B	64	LYS
1	B	80	LEU
1	B	90	VAL
1	B	91	GLN
1	B	111	VAL
1	B	116	PHE
1	B	120	LEU
1	B	138	GLU
1	B	173	LYS
1	B	205	LEU
1	B	209	LEU
1	B	214	LEU
1	B	248	GLU
1	B	250	ASP
1	B	253	THR
1	B	260	LEU
1	B	265	ASN
1	B	270	ILE
1	B	279	LEU
1	B	286	THR
1	B	293	ILE
1	B	295	LEU
1	B	297	GLU
1	B	300	GLU
1	B	301	LEU
1	B	310	LEU
1	B	347	LYS
1	B	349	LEU
1	B	377	THR
1	B	394	GLN
1	B	397	THR
1	B	422	LEU
1	B	425	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	91	GLN
1	A	103	ASN
1	A	147	ASN

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Mol	Chain	Res	Type
1	A	175	ASN
1	A	198	HIS
1	A	258	GLN
1	A	330	GLN
1	A	507	GLN
1	A	509	GLN
1	B	147	ASN
1	B	161	GLN
1	B	175	ASN
1	B	208	HIS
1	B	255	ASN
1	B	258	GLN
1	B	306	ASN
1	B	336	GLN
1	B	367	GLN
1	B	373	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](#)

1 ligand 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
2	5MA	A	561	-	33,35,35	1.63	7 (21%)	30,51,51	1.79	6 (20%)

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
2	5MA	A	561	-	-	0/8/10/10	0/5/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	561	5MA	C14-N	-2.52	1.44	1.47
2	A	561	5MA	C11-CL2	-2.24	1.69	1.74
2	A	561	5MA	C2-C1	2.01	1.40	1.36
2	A	561	5MA	C22-C23	2.87	1.43	1.36
2	A	561	5MA	C14-C15	2.96	1.55	1.51
2	A	561	5MA	C5-C4	3.40	1.47	1.42
2	A	561	5MA	C21-N20	3.78	1.40	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	561	5MA	O-C5-C	-4.89	115.20	120.25
2	A	561	5MA	C6-C11-CL2	-2.26	116.34	119.14
2	A	561	5MA	C21-N20-C18	-2.26	113.52	116.44
2	A	561	5MA	C23-C22-C21	-2.06	115.95	118.88
2	A	561	5MA	C10-C11-C6	2.36	124.84	121.69
2	A	561	5MA	O-C5-C4	6.01	125.84	117.93

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	561	5MA	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	554/563 (98%)	-0.03	18 (3%) 51 51	27, 50, 76, 99	0
1	B	400/563 (71%)	0.22	29 (7%) 18 18	28, 49, 86, 107	0
All	All	954/1126 (84%)	0.08	47 (4%) 33 34	27, 50, 82, 107	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	361	HIS	6.1
1	B	14	PRO	5.8
1	B	15	GLY	5.8
1	B	362	THR	5.4
1	B	69	THR	4.4
1	B	301	LEU	4.2
1	A	69	THR	3.9
1	B	298	GLU	3.9
1	A	284	ARG	3.8
1	B	297	GLU	3.4
1	B	292	VAL	3.4
1	A	359	GLY	3.3
1	B	70	LYS	3.3
1	B	251	SER	3.3
1	A	195	ILE	3.2
1	A	402	TRP	3.1
1	B	295	LEU	3.0
1	B	346	PHE	3.0
1	A	356	ARG	2.9
1	A	286	THR	2.9
1	B	24	TRP	2.9
1	B	278	GLN	2.9
1	B	13	LYS	2.9
1	A	70	LYS	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	240	THR	2.8
1	B	250	ASP	2.8
1	B	16	MET	2.7
1	A	221	HIS	2.7
1	B	246	LEU	2.6
1	A	52	PRO	2.6
1	A	358	ARG	2.6
1	B	232	TYR	2.6
1	B	252	TRP	2.6
1	A	360	ALA	2.6
1	B	231	GLY	2.4
1	B	284	ARG	2.3
1	B	304	ALA	2.2
1	A	137	ASN	2.2
1	A	193	LEU	2.1
1	B	279	LEU	2.1
1	B	248	GLU	2.1
1	B	294	PRO	2.1
1	A	426	TRP	2.1
1	A	15	GLY	2.1
1	B	296	THR	2.1
1	A	174	GLN	2.0
1	A	334	GLN	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(\AA^2)	Q<0.9
2	5MA	A	561	31/31	0.97	0.11	-0.82	30,41,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.