



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:29 PM GMT

PDB ID : 1KHW
Title : Crystal Structure of Rabbit Hemorrhagic Disease Virus RNA-dependent RNA polymerase complexed with Mn²⁺
Authors : Ng, K.K.; Cherney, M.M.; Vazquez, A.L.; Machin, A.; Alonso, J.M.; Parra, F.; James, M.N.
Deposited on : 2001-12-01
Resolution : 2.70 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

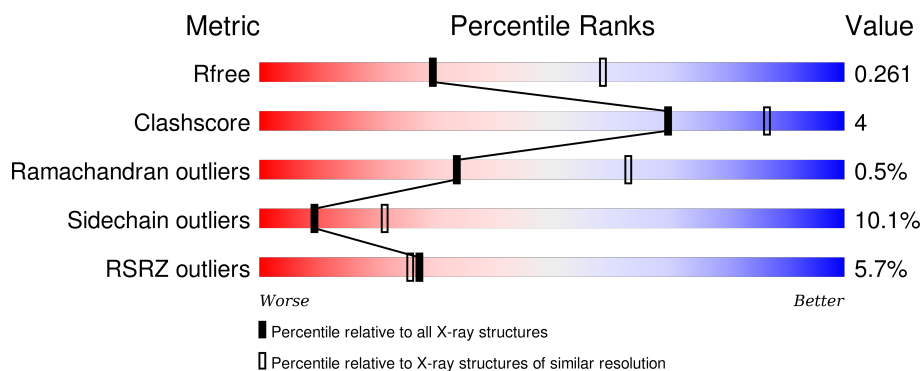
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

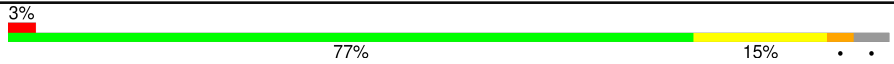

The reported resolution of this entry is 2.70 Å.

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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	516	
1	B	516	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7824 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 RNA-DIRECTED RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			3871	2464	660	719	28			
1	B	497	Total	C	N	O	S	0	0	0
			3905	2486	666	725	28			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mn	0	0
			2	2		
2	A	2	Total	Mn	0	0
			2	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total	O	0	0
			25	25		
3	B	19	Total	O	0	0
			19	19		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.69Å 119.84Å 159.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.16 – 2.70 30.44 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (31.16-2.70) 100.0 (30.44-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.238 , 0.267 0.228 , 0.261	Depositor DCC
R_{free} test set	3795 reflections (11.11%)	DCC
Wilson B-factor (Å ²)	58.0	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 37960 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7824	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.46	0/3954	0.86	13/5356 (0.2%)
1	B	0.42	0/3989	0.84	21/5403 (0.4%)
All	All	0.44	0/7943	0.85	34/10759 (0.3%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	94	ASP	CB-CG-OD2	6.35	124.01	118.30
1	B	194	ASP	CB-CG-OD2	6.18	123.86	118.30
1	B	377	ASP	CB-CG-OD2	5.91	123.62	118.30
1	A	250	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	121	ASP	CB-CG-OD2	5.79	123.52	118.30
1	B	386	ASP	CB-CG-OD2	5.62	123.36	118.30
1	A	346	ASP	CB-CG-OD2	5.57	123.31	118.30
1	B	226	ASP	CB-CG-OD2	5.53	123.28	118.30
1	B	490	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	346	ASP	CB-CG-OD2	5.47	123.23	118.30
1	A	392	ASP	CB-CG-OD2	5.39	123.15	118.30
1	A	418	ASP	CB-CG-OD2	5.37	123.14	118.30
1	B	231	ASP	CB-CG-OD2	5.36	123.13	118.30
1	B	174	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	140	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	59	ASP	CB-CG-OD2	5.32	123.09	118.30
1	B	418	ASP	CB-CG-OD2	5.30	123.07	118.30
1	B	59	ASP	CB-CG-OD2	5.29	123.06	118.30
1	B	443	ASP	CB-CG-OD2	5.24	123.02	118.30
1	B	268	ASP	CB-CG-OD2	5.24	123.01	118.30
1	A	355	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	82	ASP	CB-CG-OD2	5.21	122.98	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	ASP	CB-CG-OD2	5.18	122.97	118.30
1	B	42	ASP	CB-CG-OD2	5.17	122.95	118.30
1	A	498	ASP	CB-CG-OD2	5.16	122.95	118.30
1	A	410	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	272	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	490	ASP	CB-CG-OD2	5.12	122.90	118.30
1	B	57	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	245	ASP	CB-CG-OD2	5.08	122.88	118.30
1	B	180	ASP	CB-CG-OD2	5.07	122.86	118.30
1	A	194	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	140	ASP	CB-CG-OD2	5.04	122.83	118.30
1	B	61	ASP	CB-CG-OD2	5.04	122.83	118.30

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	3871	0	3871	40	0
1	B	3905	0	3915	30	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	25	0	0	0	0
3	B	19	0	0	0	0
All	All	7824	0	7786	70	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 (70) 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:5:PHE:HE1	1:A:167:ILE:HD11	1.18	1.04
1:A:5:PHE:CE1	1:A:167:ILE:CD1	2.49	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:PHE:HE1	1:A:167:ILE:CD1	1.82	0.91
1:A:5:PHE:CE1	1:A:167:ILE:HD11	2.06	0.86
1:A:5:PHE:CE1	1:A:167:ILE:HD12	2.16	0.81
1:A:122:LEU:HA	1:A:134:LYS:HD2	1.64	0.79
1:B:16:THR:H	1:B:76:GLN:HE22	1.34	0.72
1:A:425:GLN:O	1:A:449:ARG:NH2	2.26	0.69
1:B:49:ARG:NH1	1:B:50:PRO:O	2.29	0.66
1:B:122:LEU:HA	1:B:134:LYS:HD2	1.79	0.64
1:B:275:GLU:HG2	1:B:277:THR:HG23	1.81	0.61
1:B:108:GLN:HG3	1:B:207:ASN:HB3	1.82	0.61
1:B:49:ARG:HG3	1:B:49:ARG:HH11	1.68	0.58
1:B:425:GLN:O	1:B:449:ARG:NH2	2.37	0.58
1:B:84:VAL:HG13	1:B:88:ILE:HB	1.86	0.58
1:A:12:TYR:OH	1:A:66:ASN:ND2	2.38	0.55
1:A:67:ILE:HG21	1:A:187:LYS:HE2	1.89	0.55
1:B:108:GLN:OE1	1:B:211:LYS:NZ	2.41	0.54
1:A:424:ARG:HH11	1:A:424:ARG:HG3	1.72	0.54
1:A:386:ASP:O	1:A:388:THR:N	2.32	0.52
1:B:126:CYS:HB3	1:B:130:VAL:HB	1.92	0.51
1:B:282:SER:O	1:B:286:THR:HG23	2.11	0.51
1:A:15:ILE:HG12	1:A:69:VAL:HG13	1.93	0.50
1:A:165:HIS:HB3	1:A:167:ILE:HG13	1.92	0.50
1:A:33:ARG:HH21	1:A:174:ASP:HB3	1.78	0.49
1:A:5:PHE:CZ	1:A:167:ILE:HD12	2.47	0.49
1:A:170:CYS:HB3	1:A:295:LEU:HA	1.95	0.49
1:A:111:LEU:HB2	1:A:203:ALA:HA	1.95	0.48
1:B:49:ARG:HG3	1:B:49:ARG:NH1	2.27	0.48
1:A:258:MET:HE1	1:A:324:LEU:HD11	1.96	0.48
1:B:24:GLU:CD	1:B:24:GLU:H	2.18	0.47
1:B:153:LYS:O	1:B:157:VAL:HG12	2.15	0.47
1:B:168:TYR:HB2	1:B:293:THR:HG23	1.96	0.47
1:A:163:ALA:HB2	1:A:285:LEU:HB3	1.97	0.47
1:A:55:ARG:HG2	1:A:437:VAL:HG13	1.95	0.47
1:A:431:THR:HG22	1:A:433:SER:H	1.80	0.47
1:A:48:TYR:CD2	1:A:424:ARG:HD3	2.51	0.46
1:B:177:ARG:NH1	1:B:186:LYS:O	2.49	0.46
1:B:61:ASP:N	1:B:61:ASP:OD1	2.47	0.46
1:B:411:ILE:HG12	1:B:411:ILE:O	2.14	0.46
1:B:343:LEU:HA	1:B:343:LEU:HD22	1.81	0.46
1:B:122:LEU:HB3	1:B:135:ILE:HB	1.98	0.45
1:A:258:MET:HE2	1:A:259:SER:H	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:472:CYS:HA	1:B:475:LEU:HB2	1.97	0.45
1:A:225:VAL:HG21	1:A:235:ILE:HD11	1.99	0.45
1:A:195:VAL:O	1:A:199:VAL:HG23	2.17	0.45
1:A:264:ARG:HG2	1:A:288:LYS:HG2	1.99	0.45
1:B:396:LEU:HD13	1:B:413:TRP:HH2	1.82	0.44
1:A:56:SER:HB2	1:A:437:VAL:HG12	2.00	0.44
1:A:301:GLN:H	1:A:301:GLN:CD	2.21	0.44
1:B:49:ARG:HH11	1:B:49:ARG:CG	2.28	0.43
1:B:231:ASP:O	1:B:235:ILE:HG23	2.19	0.43
1:A:36:LYS:NZ	1:A:427:GLU:OE2	2.52	0.43
1:B:387:LYS:HE3	1:B:387:LYS:HB2	1.83	0.43
1:A:431:THR:HG22	1:A:432:THR:N	2.34	0.43
1:B:100:ARG:NH2	1:B:272:ASP:OD2	2.52	0.42
1:A:317:ASN:HB3	1:A:353:GLY:HA2	2.01	0.42
1:B:37:VAL:HA	1:B:38:PRO:HD3	1.92	0.42
1:B:247:LEU:HB3	1:B:358:TYR:HD1	1.84	0.42
1:A:5:PHE:HB3	1:A:21:VAL:O	2.20	0.42
1:A:118:ASN:HD22	1:A:118:ASN:HA	1.70	0.41
1:A:84:VAL:HG13	1:A:88:ILE:HB	2.03	0.41
1:A:385:ALA:O	1:A:387:LYS:N	2.53	0.41
1:B:264:ARG:HA	1:B:288:LYS:HG2	2.01	0.41
1:A:295:LEU:HB3	1:A:298:MET:HG2	2.01	0.41
1:A:10:ILE:HD12	1:A:298:MET:HB2	2.02	0.41
1:B:222:ALA:O	1:B:225:VAL:HB	2.20	0.41
1:A:228:THR:HG23	1:A:499:ARG:HB3	2.03	0.41
1:A:494:LYS:HE3	1:A:494:LYS:HB3	1.87	0.41
1:A:313:THR:O	1:A:317:ASN:HB2	2.21	0.41

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	489/516 (95%)	461 (94%)	25 (5%)	3 (1%)	30	59
1	B	495/516 (96%)	477 (96%)	16 (3%)	2 (0%)	39	69
All	All	984/1032 (95%)	938 (95%)	41 (4%)	5 (0%)	34	63

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	387	LYS
1	B	21	VAL
1	B	296	ASP
1	A	62	LYS
1	A	139	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	423/446 (95%)	385 (91%)	38 (9%)	12	27
1	B	427/446 (96%)	379 (89%)	48 (11%)	7	17
All	All	850/892 (95%)	764 (90%)	86 (10%)	9	21

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	PHE
1	A	30	SER
1	A	52	ASN
1	A	55	ARG
1	A	62	LYS
1	A	84	VAL
1	A	102	LEU
1	A	111	LEU
1	A	120	LEU
1	A	121	ASP
1	A	125	SER

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Mol	Chain	Res	Type
1	A	138	VAL
1	A	147	LEU
1	A	173	LYS
1	A	194	ASP
1	A	215	VAL
1	A	228	THR
1	A	234	VAL
1	A	242	LYS
1	A	247	LEU
1	A	285	LEU
1	A	298	MET
1	A	301	GLN
1	A	304	ARG
1	A	306	LEU
1	A	351	THR
1	A	352	TYR
1	A	358	TYR
1	A	361	THR
1	A	387	LYS
1	A	400	SER
1	A	418	ASP
1	A	432	THR
1	A	433	SER
1	A	437	VAL
1	A	449	ARG
1	A	451	VAL
1	A	493	ARG
1	B	5	PHE
1	B	6	CYS
1	B	8	GLU
1	B	15	ILE
1	B	24	GLU
1	B	36	LYS
1	B	42	ASP
1	B	49	ARG
1	B	61	ASP
1	B	64	LEU
1	B	81	LEU
1	B	84	VAL
1	B	108	GLN
1	B	139	LYS
1	B	147	LEU

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Mol	Chain	Res	Type
1	B	157	VAL
1	B	225	VAL
1	B	236	ILE
1	B	247	LEU
1	B	252	SER
1	B	279	LEU
1	B	293	THR
1	B	295	LEU
1	B	302	THR
1	B	303	LYS
1	B	326	SER
1	B	335	GLU
1	B	343	LEU
1	B	352	TYR
1	B	358	TYR
1	B	368	LEU
1	B	373	GLU
1	B	383	THR
1	B	386	ASP
1	B	388	THR
1	B	391	ILE
1	B	399	ILE
1	B	400	SER
1	B	403	LYS
1	B	410	ASP
1	B	411	ILE
1	B	418	ASP
1	B	437	VAL
1	B	449	ARG
1	B	477	ARG
1	B	494	LYS
1	B	496	LEU
1	B	500	LYS

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	18	HIS
1	A	52	ASN
1	A	66	ASN
1	A	118	ASN
1	A	479	GLN

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Mol	Chain	Res	Type
1	B	52	ASN
1	B	73	GLN
1	B	76	GLN
1	B	118	ASN
1	B	150	HIS
1	B	237	ASN
1	B	342	ASN
1	B	374	ASN
1	B	465	GLN
1	B	479	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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

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	493/516 (95%)	0.07	16 (3%) 51 51	12, 26, 48, 64	0
1	B	497/516 (96%)	0.38	40 (8%) 15 13	13, 26, 48, 61	0
All	All	990/1032 (95%)	0.22	56 (5%) 27 26	12, 26, 48, 64	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	VAL	7.4
1	B	385	ALA	7.4
1	B	386	ASP	7.3
1	A	388	THR	6.4
1	B	390	PHE	5.6
1	B	389	GLU	5.0
1	A	179	LEU	4.8
1	B	140	ASP	4.8
1	B	388	THR	4.6
1	B	122	LEU	4.6
1	A	386	ASP	4.5
1	B	396	LEU	4.3
1	B	411	ILE	4.3
1	B	5	PHE	4.1
1	B	334	ALA	4.0
1	A	55	ARG	3.7
1	B	141	GLY	3.4
1	B	387	LYS	3.3
1	A	385	ALA	3.2
1	B	7	GLY	3.2
1	B	391	ILE	3.2
1	B	23	ALA	3.2
1	B	338	LEU	3.1
1	A	389	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	6	CYS	3.1
1	A	138	VAL	3.0
1	B	397	ASN	3.0
1	B	104	LYS	2.9
1	B	246	PHE	2.9
1	A	411	ILE	2.7
1	B	86	GLU	2.7
1	B	366	SER	2.7
1	B	399	ILE	2.7
1	B	383	THR	2.7
1	B	22	GLY	2.7
1	B	336	ILE	2.6
1	B	249	LEU	2.6
1	B	103	THR	2.6
1	A	49	ARG	2.6
1	B	210	TYR	2.5
1	A	135	ILE	2.4
1	A	432	THR	2.4
1	B	342	ASN	2.3
1	B	107	ARG	2.3
1	B	101	PHE	2.3
1	B	320	CYS	2.3
1	B	139	LYS	2.3
1	B	384	ALA	2.2
1	A	112	ASN	2.2
1	A	64	LEU	2.2
1	B	393	VAL	2.1
1	A	5	PHE	2.1
1	A	501	ARG	2.1
1	A	8	GLU	2.0
1	B	214	MET	2.0
1	B	341	SER	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MN	B	604	1/1	0.76	0.10	-	55,55,55,55	1
2	MN	B	603	1/1	0.90	0.75	-	84,84,84,84	1
2	MN	A	601	1/1	0.94	0.09	-	64,64,64,64	0
2	MN	A	602	1/1	0.80	0.09	-	69,69,69,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.