



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:37 PM UTC

PDB ID : 2V83 / pdb_00002v83
Title : Crystal structure of RAG2-PHD finger in complex with H3K4me3 peptide
Authors : Ramon-Maiques, S.; Yang, W.
Deposited on : 2007-08-02
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

<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

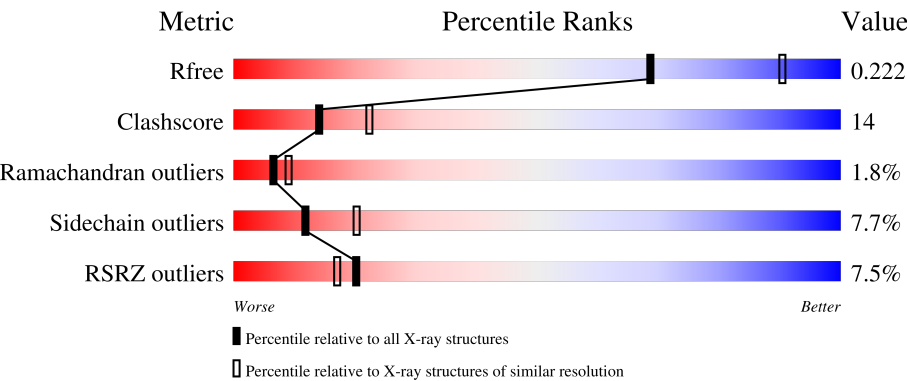
1 Overall quality at a glance i

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	82	% 67% 18% • 12%
1	B	82	% 70% 24% • •
1	C	82	20% 39% 37% 10% • 12%
2	D	9	44% 44% 11%
2	E	9	78% 22%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2177 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 VDJ RECOMBINATION-ACTIVATING PROTEIN 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	72	Total	C	N	O	S	0	0	1
			575	362	96	109	8			
1	B	79	Total	C	N	O	S	0	0	0
			635	401	106	120	8			
1	C	72	Total	C	N	O	S	0	0	1
			579	368	95	108	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	406	GLY	-	expression tag	UNP P21784
A	407	PRO	-	expression tag	UNP P21784
A	408	LEU	-	expression tag	UNP P21784
A	409	GLY	-	expression tag	UNP P21784
A	410	SER	-	expression tag	UNP P21784
A	411	PRO	-	expression tag	UNP P21784
A	412	GLU	-	expression tag	UNP P21784
A	413	PHE	-	expression tag	UNP P21784
B	406	GLY	-	expression tag	UNP P21784
B	407	PRO	-	expression tag	UNP P21784
B	408	LEU	-	expression tag	UNP P21784
B	409	GLY	-	expression tag	UNP P21784
B	410	SER	-	expression tag	UNP P21784
B	411	PRO	-	expression tag	UNP P21784
B	412	GLU	-	expression tag	UNP P21784
B	413	PHE	-	expression tag	UNP P21784
C	406	GLY	-	expression tag	UNP P21784
C	407	PRO	-	expression tag	UNP P21784
C	408	LEU	-	expression tag	UNP P21784
C	409	GLY	-	expression tag	UNP P21784
C	410	SER	-	expression tag	UNP P21784
C	411	PRO	-	expression tag	UNP P21784
C	412	GLU	-	expression tag	UNP P21784

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	413	PHE	-	expression tag	UNP P21784

- Molecule 2 is a protein called HISTONE H3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	9	Total	C	N	O	0	0	0
			77	46	18	13			
2	E	9	Total	C	N	O	0	0	0
			71	43	15	13			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		
3	C	2	Total	Zn	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	81	Total	O	0	0
			81	81		
4	B	91	Total	O	0	0
			91	91		
4	C	48	Total	O	0	0
			48	48		
4	D	8	Total	O	0	0
			8	8		
4	E	6	Total	O	0	0
			6	6		

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.

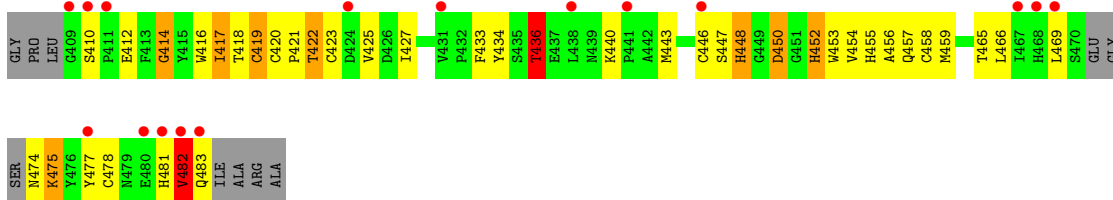
• Molecule 1: VDJ RECOMBINATION-ACTIVATING PROTEIN 2



• Molecule 1: VDJ RECOMBINATION-ACTIVATING PROTEIN 2



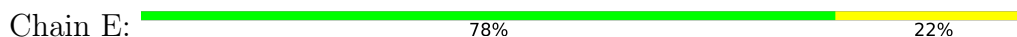
• Molecule 1: VDJ RECOMBINATION-ACTIVATING PROTEIN 2



• Molecule 2: HISTONE H3



• Molecule 2: HISTONE H3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.78Å 46.84Å 56.96Å 90.00° 101.46° 90.00°	Depositor
Resolution (Å)	30.00 – 2.40 30.00 – 2.41	Depositor EDS
% Data completeness (in resolution range)	99.4 (30.00-2.40) 91.2 (30.00-2.41)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.42Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.230 0.184 , 0.222	Depositor DCC
R_{free} test set	530 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	16.9	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.053 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2177	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.04% 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: M3L, ZN

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	0.47	1/594 (0.2%)	0.93	3/812 (0.4%)
1	B	0.48	0/656	0.95	3/894 (0.3%)
1	C	0.53	1/599 (0.2%)	1.07	7/818 (0.9%)
2	D	0.52	0/64	1.30	0/82
2	E	0.47	0/58	0.77	0/75
All	All	0.49	2/1971 (0.1%)	0.99	13/2681 (0.5%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	482	VAL	C-N	-5.65	1.25	1.33
1	A	484	ILE	C-N	-5.64	1.25	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	450	ASP	N-CA-C	-7.54	103.00	111.07
1	B	425	VAL	N-CA-C	7.49	119.11	109.30
1	B	414	GLY	N-CA-C	6.87	123.26	115.08
1	C	425	VAL	N-CA-C	6.73	118.51	108.54
1	C	412	GLU	N-CA-C	5.94	117.75	111.28
1	C	481	HIS	N-CA-C	5.86	121.15	113.30
1	C	436	THR	N-CA-CB	-5.59	101.04	110.49
1	C	419	CYS	N-CA-C	5.57	118.11	111.71
1	A	419	CYS	N-CA-C	5.56	119.32	112.54
1	A	430	TRP	N-CA-C	5.50	118.21	110.23
1	C	414	GLY	N-CA-C	5.36	119.67	112.17
1	B	419	CYS	N-CA-C	5.09	117.73	111.82
1	A	418	THR	N-CA-C	-5.03	99.13	107.99

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	575	0	510	9	0
1	B	635	0	569	15	0
1	C	579	0	515	26	0
2	D	77	0	91	5	0
2	E	71	0	80	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
4	A	81	0	0	1	0
4	B	91	0	0	4	0
4	C	48	0	0	1	0
4	D	8	0	0	1	0
4	E	6	0	0	0	0
All	All	2177	0	1765	53	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 14.

All (53) 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:B:486:ARG:HG2	1:B:486:ARG:O	1.77	0.84
1:C:434:TYR:HE1	1:C:456:ALA:HB3	1.56	0.71
1:C:418:THR:HA	1:C:423:CYS:SG	2.33	0.69
2:D:2:ARG:HD2	4:D:2001:HOH:O	1.95	0.65
1:C:448:HIS:C	1:C:448:HIS:HD1	2.05	0.64
1:B:420:CYS:HB2	1:B:421:PRO:HD2	1.79	0.63
1:B:419:CYS:HB2	1:B:454:VAL:HA	1.80	0.62
1:B:426:ASP:OD2	1:B:428:ASN:HB2	2.03	0.59
1:C:448:HIS:C	1:C:448:HIS:ND1	2.61	0.58
1:C:448:HIS:ND1	1:C:448:HIS:O	2.37	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:ARG:O	1:B:486:ARG:CG	2.51	0.56
1:C:427:ILE:O	2:D:4:M3L:HE3	2.06	0.56
1:A:419:CYS:HB2	1:A:454:VAL:HA	1.88	0.56
2:E:9:LYS:OXT	2:E:9:LYS:HG2	2.06	0.53
1:A:425:VAL:HG21	1:A:455:HIS:CG	2.43	0.53
1:C:443:MET:HE2	1:C:453:TRP:HB3	1.90	0.53
1:C:419:CYS:HB2	1:C:454:VAL:HA	1.91	0.52
1:C:459:MET:SD	1:C:478:CYS:HB2	2.50	0.52
4:B:2049:HOH:O	1:C:414:GLY:HA2	2.11	0.50
1:B:486:ARG:HB2	4:B:2081:HOH:O	2.10	0.50
1:B:439:ASN:HD22	1:B:439:ASN:H	1.60	0.50
1:A:442:ALA:HA	2:D:5:GLN:HB2	1.95	0.49
1:C:420:CYS:HB2	1:C:421:PRO:CD	2.42	0.49
1:C:482:VAL:HG23	1:C:483:GLN:N	2.28	0.49
1:C:446:CYS:SG	1:C:477:TYR:HA	2.54	0.48
1:C:465:THR:O	1:C:469:LEU:HG	2.15	0.47
1:B:484:ILE:O	1:B:485:ALA:C	2.57	0.47
1:C:456:ALA:HB1	1:C:466:LEU:HD12	1.97	0.47
1:B:440:LYS:NZ	4:B:2041:HOH:O	2.46	0.46
1:A:483:GLN:HA	4:A:2077:HOH:O	2.15	0.46
1:A:443:MET:HE3	1:A:455:HIS:CD2	2.51	0.46
1:B:420:CYS:H	1:B:423:CYS:HB2	1.80	0.46
1:B:421:PRO:HG2	4:B:2018:HOH:O	2.17	0.45
1:C:446:CYS:HB3	1:C:452:HIS:CD2	2.52	0.44
1:B:446:CYS:HB3	1:B:452:HIS:CD2	2.53	0.44
1:C:456:ALA:CB	1:C:466:LEU:HD12	2.49	0.43
1:B:482:VAL:HG13	1:B:483:GLN:HG3	2.00	0.42
1:C:434:TYR:CE1	1:C:456:ALA:HB3	2.45	0.42
1:C:455:HIS:HB2	1:C:458:CYS:SG	2.60	0.42
2:D:8:ARG:CD	2:D:8:ARG:H	2.26	0.42
1:A:420:CYS:HB2	1:A:480:GLU:OE2	2.20	0.42
1:A:425:VAL:HB	1:A:455:HIS:CE1	2.54	0.41
1:A:478:CYS:O	1:A:482:VAL:HG22	2.21	0.41
1:C:422:THR:O	1:C:423:CYS:C	2.63	0.41
1:A:463:GLU:O	1:A:467:ILE:HG12	2.21	0.41
1:C:433:PHE:CD2	1:C:457:GLN:HG3	2.56	0.41
1:C:459:MET:HE1	4:C:2047:HOH:O	2.20	0.41
1:C:474:ASN:HB2	1:C:475:LYS:H	1.67	0.41
1:C:417:ILE:HD11	1:C:443:MET:HE1	2.03	0.41
1:C:452:HIS:CD2	1:C:452:HIS:N	2.89	0.41
1:C:416:TRP:HB2	2:D:2:ARG:NH2	2.37	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:TRP:CG	2:E:4:M3L:HM33	2.57	0.40
1:B:439:ASN:HD22	1:B:439:ASN:C	2.28	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	70/82 (85%)	66 (94%)	4 (6%)	0	100	100
1	B	77/82 (94%)	70 (91%)	6 (8%)	1 (1%)	9	14
1	C	68/82 (83%)	57 (84%)	8 (12%)	3 (4%)	2	1
2	D	6/9 (67%)	6 (100%)	0	0	100	100
2	E	6/9 (67%)	6 (100%)	0	0	100	100
All	All	227/264 (86%)	205 (90%)	18 (8%)	4 (2%)	6	9

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	482	VAL
1	B	410	SER
1	C	410	SER
1	C	436	THR

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	64/72 (89%)	62 (97%)	2 (3%)	35	57
1	B	69/72 (96%)	66 (96%)	3 (4%)	26	44
1	C	65/72 (90%)	56 (86%)	9 (14%)	3	4
2	D	6/6 (100%)	4 (67%)	2 (33%)	0	0
2	E	5/6 (83%)	5 (100%)	0	100	100
All	All	209/228 (92%)	193 (92%)	16 (8%)	12	20

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	466	LEU
1	A	482	VAL
1	B	439	ASN
1	B	464	ARG
1	B	466	LEU
1	C	417	ILE
1	C	422	THR
1	C	436	THR
1	C	440	LYS
1	C	447	SER
1	C	448	HIS
1	C	450	ASP
1	C	452	HIS
1	C	475	LYS
2	D	8	ARG
2	D	9	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	439	ASN
1	A	448	HIS
1	A	474	ASN
1	A	483	GLN
1	B	439	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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	M3L	D	4	2	10,11,12	0.46	0	9,14,16	0.42	0
2	M3L	E	4	2	10,11,12	0.43	0	9,14,16	0.65	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
2	M3L	D	4	2	-	1/9/10/12	-
2	M3L	E	4	2	-	0/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	4	M3L	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	M3L	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	4	M3L	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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, 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	72/82 (87%)	-0.42	1 (1%) 73 69	6, 13, 28, 58	0
1	B	79/82 (96%)	-0.22	1 (1%) 75 71	7, 17, 32, 45	0
1	C	72/82 (87%)	1.51	16 (22%) 2 2	40, 57, 82, 84	0
2	D	8/9 (88%)	0.16	0 100 100	12, 20, 51, 54	0
2	E	8/9 (88%)	0.13	0 100 100	14, 20, 41, 50	0
All	All	239/264 (90%)	0.27	18 (7%) 20 17	6, 21, 76, 84	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	483	GLN	5.8
1	A	485	ALA	5.1
1	C	438	LEU	4.3
1	B	487	ALA	3.8
1	C	482	VAL	3.7
1	C	409	GLY	3.5
1	C	467	ILE	3.0
1	C	410	SER	2.9
1	C	477	TYR	2.8
1	C	480	GLU	2.8
1	C	468	HIS	2.5
1	C	411	PRO	2.4
1	C	431	VAL	2.4
1	C	481	HIS	2.3
1	C	441	PRO	2.2
1	C	446	CYS	2.2
1	C	424	ASP	2.2
1	C	469	LEU	2.1

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
2	M3L	E	4	12/13	0.93	0.09	15,16,22,24	0
2	M3L	D	4	12/13	0.95	0.08	9,11,16,17	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
3	ZN	A	1486	1/1	0.99	0.03	12,12,12,12	0
3	ZN	B	1489	1/1	0.99	0.02	16,16,16,16	0
3	ZN	C	1484	1/1	0.99	0.03	41,41,41,41	0
3	ZN	C	1485	1/1	0.99	0.04	57,57,57,57	0
3	ZN	B	1488	1/1	1.00	0.02	15,15,15,15	0
3	ZN	A	1487	1/1	1.00	0.03	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.