



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 11:16 PM UTC

PDB ID : 7B1F / pdb_00007b1f
Title : Orthorhombic P212121 Structure of Human Mad1 C-terminal Domain in Complex with Phosphorylated Bub1 CD1 Domain
Authors : Fischer, E.; Bellini, D.; Barford, D.
Deposited on : 2020-11-24
Resolution : 1.75 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

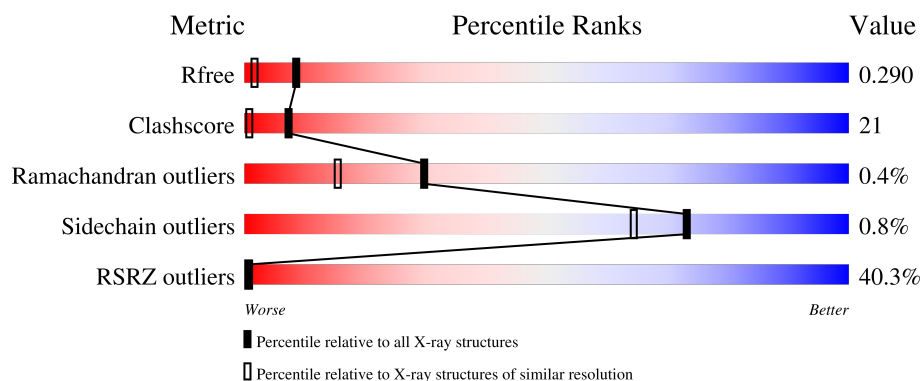
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	122	19% 84% 16%
1	B	122	46% 80% 16%
2	C	26	65% 38% 31% 12% 19%
2	D	26	62% 42% 27% 12% 19%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SEP	C	459	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2408 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 Mitotic spindle assembly checkpoint protein MAD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	118	Total	C	N	O	S	0	0	0
			957	606	162	186	3			
1	A	122	Total	C	N	O	S	0	0	0
			978	618	166	191	3			

- Molecule 2 is a protein called Mitotic checkpoint serine/threonine-protein kinase BUB1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	21	Total	C	N	O	P S	0	0	0
			167	103	26	34	2 2			
2	D	21	Total	C	N	O	P S	0	0	0
			167	103	26	34	2 2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	480	SER	-	expression tag	UNP O43683
D	480	SER	-	expression tag	UNP O43683

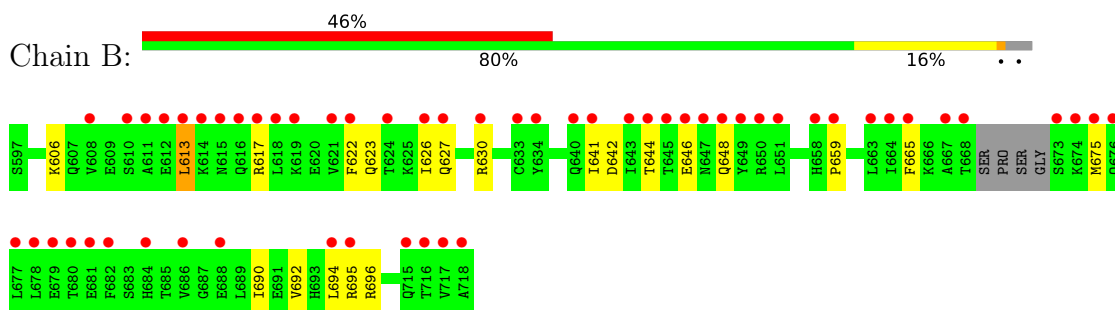
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	46	Total	O	0	0
			46	46		
3	A	91	Total	O	0	0
			91	91		
3	C	1	Total	O	0	0
			1	1		
3	D	1	Total	O	0	0
			1	1		

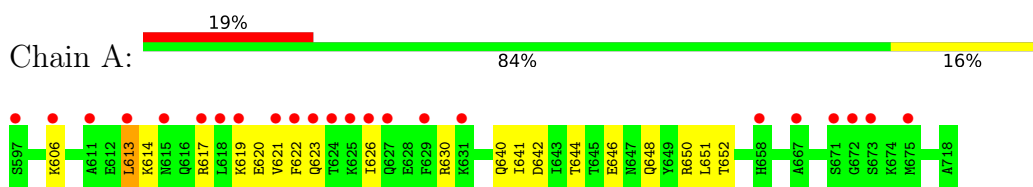
3 Residue-property plots [i](#)

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.

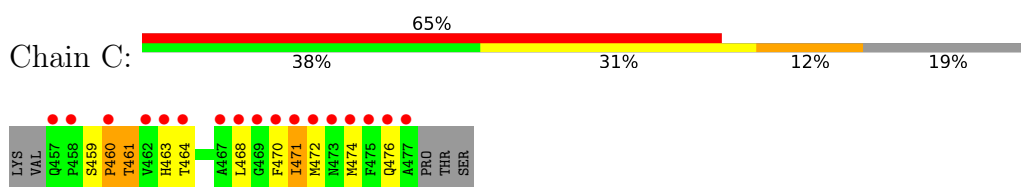
- Molecule 1: Mitotic spindle assembly checkpoint protein MAD1



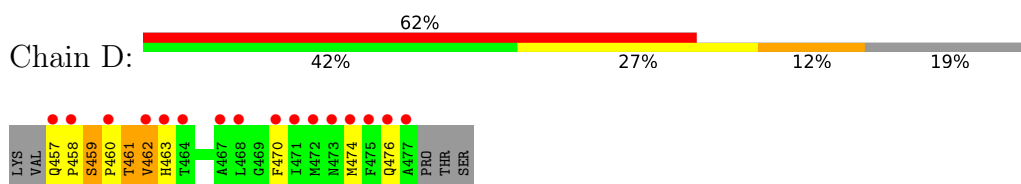
- Molecule 1: Mitotic spindle assembly checkpoint protein MAD1



- Molecule 2: Mitotic checkpoint serine/threonine-protein kinase BUB1



- Molecule 2: Mitotic checkpoint serine/threonine-protein kinase BUB1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	34.22Å 80.49Å 134.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.25 – 1.75 40.25 – 1.75	Depositor EDS
% Data completeness (in resolution range)	85.8 (40.25-1.75) 85.7 (40.25-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.83 (at 1.75Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.256 , 0.287 0.258 , 0.290	Depositor DCC
R_{free} test set	1615 reflections (4.21%)	wwPDB-VP
Wilson B-factor (Å ²)	37.4	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 30.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2408	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.21% 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: SEP, TPO

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	0.58	3/993 (0.3%)	0.61	1/1337 (0.1%)
1	B	0.39	0/970	0.68	0/1304
2	C	0.98	0/148	1.35	2/195 (1.0%)
2	D	1.08	0/148	1.09	1/195 (0.5%)
All	All	0.59	3/2259 (0.1%)	0.74	4/3031 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	613	LEU	CA-CB	-5.95	1.45	1.54
1	A	613	LEU	C-O	-5.78	1.16	1.23
1	A	613	LEU	N-CA	-5.44	1.39	1.45

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	614	LYS	N-CA-C	-12.13	98.14	111.36
2	C	470	PHE	N-CA-C	-7.95	103.19	113.12
2	D	462	VAL	CB-CA-C	-7.37	97.40	111.40
2	C	471	ILE	N-CA-C	5.53	115.98	110.23

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	978	0	984	38	0
1	B	957	0	967	32	2
2	C	167	0	151	46	0
2	D	167	0	151	21	0
3	A	91	0	0	0	0
3	B	46	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
All	All	2408	0	2253	93	2

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 21.

All (93) 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:623:GLN:HE22	2:C:474:MET:CG	1.26	1.45
1:A:622:PHE:CD1	2:C:471:ILE:HD13	1.57	1.39
1:B:623:GLN:NE2	2:C:474:MET:HG3	1.44	1.30
1:A:623:GLN:HG3	2:D:474:MET:HE1	1.19	1.18
1:B:623:GLN:NE2	2:C:474:MET:CG	1.99	1.14
1:A:623:GLN:HG3	2:D:474:MET:CE	1.76	1.14
2:C:468:LEU:HA	2:C:471:ILE:HG22	1.20	1.13
1:B:623:GLN:NE2	2:C:474:MET:SD	2.25	1.08
1:B:626:ILE:HD11	2:C:474:MET:HB2	1.30	1.07
2:C:468:LEU:HA	2:C:471:ILE:CG2	1.85	1.07
1:B:626:ILE:HD11	2:C:474:MET:CB	1.88	1.03
1:A:622:PHE:HD1	2:C:471:ILE:HD13	0.83	0.97
1:A:622:PHE:CD1	2:C:471:ILE:CD1	2.48	0.97
2:C:459:SEP:O3P	2:C:460:PRO:HD2	1.65	0.95
1:A:622:PHE:HD1	2:C:471:ILE:CD1	1.79	0.95
2:D:457:GLN:N	2:D:458:PRO:HD3	1.83	0.94
2:C:468:LEU:CA	2:C:471:ILE:HG22	1.99	0.91
2:C:459:SEP:HB2	2:C:460:PRO:CD	2.02	0.88
2:C:468:LEU:HD23	2:C:471:ILE:HG21	1.53	0.87
1:A:619:LYS:HG2	2:D:470:PHE:CE2	2.09	0.86
2:C:459:SEP:HB2	2:C:460:PRO:HD2	1.57	0.85
1:A:630:ARG:NH2	1:A:640:GLN:OE1	2.10	0.85
2:C:459:SEP:CB	2:C:460:PRO:CD	2.54	0.84
1:A:619:LYS:HG2	2:D:470:PHE:HE2	1.45	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:468:LEU:HD23	2:C:471:ILE:CG2	2.11	0.81
1:B:622:PHE:CE1	1:A:626:ILE:HD11	2.22	0.75
1:B:623:GLN:CD	2:C:474:MET:HG3	2.10	0.75
1:B:627:GLN:HE21	2:C:474:MET:CE	2.01	0.73
1:B:690:ILE:HG23	1:B:694:LEU:HD12	1.71	0.72
1:A:622:PHE:O	1:A:626:ILE:HG12	1.90	0.71
1:A:644:THR:HG23	1:A:646:GLU:H	1.56	0.71
2:C:459:SEP:O3P	2:C:460:PRO:CD	2.40	0.70
1:A:621:VAL:CG1	2:C:471:ILE:HG21	2.21	0.70
1:A:623:GLN:CG	2:D:474:MET:HE1	2.12	0.68
1:B:626:ILE:HD11	2:C:474:MET:HB3	1.75	0.67
1:A:623:GLN:HG3	2:D:474:MET:HE2	1.70	0.67
1:B:617:ARG:HG3	2:D:463:HIS:HB3	1.78	0.66
1:B:626:ILE:CD1	2:C:474:MET:HB2	2.20	0.64
2:D:459:SEP:HB2	2:D:460:PRO:HD2	1.79	0.64
2:D:457:GLN:N	2:D:458:PRO:CD	2.61	0.62
1:B:622:PHE:CZ	1:A:626:ILE:HD11	2.35	0.62
1:A:644:THR:HG22	1:A:648:GLN:H	1.64	0.62
1:B:695:ARG:NH1	1:B:696:ARG:HE	2.00	0.59
2:C:468:LEU:HD22	2:C:472:MET:HE3	1.85	0.58
1:B:623:GLN:CD	2:C:474:MET:CG	2.72	0.57
1:A:622:PHE:HA	2:C:471:ILE:CD1	2.35	0.56
2:D:461:TPO:HG22	2:D:461:TPO:O1P	2.06	0.55
1:A:630:ARG:NH2	1:A:642:ASP:OD1	2.40	0.54
1:B:622:PHE:CZ	2:C:471:ILE:HD12	2.43	0.54
2:D:462:VAL:HG12	2:D:462:VAL:O	2.08	0.53
1:B:630:ARG:HH22	1:B:642:ASP:HA	1.73	0.53
1:B:627:GLN:HE21	2:C:474:MET:HE1	1.72	0.53
1:A:650:ARG:NH2	2:D:476:GLN:OE1	2.36	0.52
1:A:621:VAL:HG11	2:C:471:ILE:HG21	1.91	0.52
2:C:461:TPO:HG22	2:C:461:TPO:O1P	2.09	0.52
2:C:468:LEU:C	2:C:471:ILE:HG22	2.35	0.52
1:A:621:VAL:HG12	2:C:471:ILE:HG12	1.91	0.52
1:A:644:THR:HG23	1:A:646:GLU:N	2.24	0.52
1:A:617:ARG:HA	1:A:620:GLU:HG2	1.92	0.51
1:B:622:PHE:CE2	2:C:471:ILE:HD12	2.47	0.50
1:B:644:THR:OG1	1:B:648:GLN:HB2	2.12	0.49
2:C:461:TPO:HG21	2:C:461:TPO:O2P	2.10	0.49
1:B:665:PHE:HB3	1:B:675:MET:HE2	1.92	0.49
1:A:619:LYS:CG	2:D:470:PHE:HE2	2.22	0.49
1:A:622:PHE:HA	2:C:471:ILE:HD13	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:640:GLN:HB3	1:A:652:THR:OG1	2.15	0.47
1:A:623:GLN:HA	2:D:474:MET:HE2	1.97	0.47
1:A:622:PHE:HA	2:C:471:ILE:HD11	1.97	0.47
1:A:623:GLN:CG	2:D:474:MET:CE	2.69	0.46
1:B:630:ARG:NH2	1:B:641:ILE:O	2.49	0.46
1:B:613:LEU:O	1:B:617:ARG:HG2	2.16	0.46
2:C:468:LEU:HA	2:C:471:ILE:HG21	1.86	0.45
1:B:622:PHE:CZ	2:C:471:ILE:CD1	2.99	0.45
1:B:692:VAL:O	1:B:696:ARG:HB2	2.16	0.45
1:B:626:ILE:CD1	2:C:474:MET:CB	2.79	0.45
2:C:461:TPO:O3P	2:C:463:HIS:HB2	2.17	0.45
2:C:459:SEP:HA	2:C:460:PRO:HD3	1.33	0.44
1:B:695:ARG:HH11	1:B:696:ARG:HE	1.64	0.44
1:A:641:ILE:HG12	1:A:651:LEU:HG	2.00	0.44
1:B:623:GLN:NE2	2:C:474:MET:CE	2.80	0.43
1:B:617:ARG:HG3	2:D:463:HIS:CB	2.47	0.43
1:A:644:THR:HG22	1:A:648:GLN:N	2.32	0.43
1:B:606:LYS:HE2	1:B:606:LYS:HB3	1.75	0.43
1:A:606:LYS:HD2	1:A:606:LYS:N	2.34	0.43
1:A:623:GLN:CG	2:D:474:MET:HE2	2.45	0.43
1:A:644:THR:HA	2:D:476:GLN:HE22	1.84	0.42
1:A:622:PHE:O	1:A:626:ILE:CG1	2.63	0.42
1:A:613:LEU:HA	1:A:613:LEU:HD12	1.72	0.42
2:D:461:TPO:O2P	2:D:461:TPO:HG21	2.20	0.41
1:A:617:ARG:HG2	2:C:464:THR:OG1	2.21	0.41
2:D:461:TPO:O3P	2:D:463:HIS:ND1	2.48	0.41
1:B:675:MET:HE2	1:B:675:MET:HB3	1.94	0.40
2:C:461:TPO:O1P	2:C:461:TPO:C	2.70	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:646:GLU:OE2	1:B:659:PRO:CG[4_455]	1.27	0.93
1:B:646:GLU:OE2	1:B:659:PRO:CD[4_455]	1.67	0.53

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	120/122 (98%)	119 (99%)	1 (1%)	0	100	100
1	B	114/122 (93%)	113 (99%)	1 (1%)	0	100	100
2	C	17/26 (65%)	15 (88%)	1 (6%)	1 (6%)	1	0
2	D	17/26 (65%)	17 (100%)	0	0	100	100
All	All	268/296 (90%)	264 (98%)	3 (1%)	1 (0%)	30	15

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	460	PRO

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	110/111 (99%)	110 (100%)	0	100	100
1	B	108/111 (97%)	107 (99%)	1 (1%)	70	60
2	C	15/21 (71%)	14 (93%)	1 (7%)	15	3
2	D	15/21 (71%)	15 (100%)	0	100	100
All	All	248/264 (94%)	246 (99%)	2 (1%)	73	64

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	613	LEU
2	C	476	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	623	GLN
1	B	627	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	TPO	D	461	2	8,10,11	2.05	2 (25%)	10,14,16	2.37	4 (40%)
2	TPO	C	461	2	8,10,11	1.59	2 (25%)	10,14,16	2.09	3 (30%)
2	SEP	C	459	2	8,9,10	0.97	0	7,12,14	1.25	0
2	SEP	D	459	2	8,9,10	1.14	0	7,12,14	1.69	2 (28%)

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	TPO	D	461	2	-	3/9/11/13	-
2	TPO	C	461	2	-	6/9/11/13	-
2	SEP	C	459	2	-	4/6/8/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SEP	D	459	2	-	3/6/8/10	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	461	TPO	P-OG1	-4.27	1.52	1.59
2	C	461	TPO	P-O3P	-2.69	1.44	1.54
2	C	461	TPO	P-O2P	-2.66	1.44	1.54
2	D	461	TPO	P-O3P	-2.56	1.45	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	461	TPO	P-OG1-CB	-5.06	109.58	123.33
2	C	461	TPO	P-OG1-CB	-4.37	111.44	123.33
2	D	459	SEP	OG-CB-CA	3.48	111.53	108.14
2	D	461	TPO	CG2-CB-CA	3.30	119.69	113.26
2	C	461	TPO	O2P-P-O1P	3.17	123.17	110.83
2	D	461	TPO	CB-CA-N	-2.55	97.50	114.17
2	D	461	TPO	O3P-P-OG1	-2.13	97.55	105.85
2	D	459	SEP	OG-P-O1P	-2.06	100.87	106.44
2	C	461	TPO	O3P-P-O1P	-2.03	102.94	110.83

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	459	SEP	CB-OG-P-O1P
2	C	459	SEP	CB-OG-P-O2P
2	C	459	SEP	CB-OG-P-O3P
2	C	461	TPO	N-CA-CB-OG1
2	C	461	TPO	O-C-CA-CB
2	C	461	TPO	CG2-CB-OG1-P
2	D	459	SEP	CB-OG-P-O1P
2	D	459	SEP	CB-OG-P-O2P
2	D	459	SEP	CB-OG-P-O3P
2	D	461	TPO	N-CA-CB-CG2
2	D	461	TPO	N-CA-CB-OG1
2	D	461	TPO	CG2-CB-OG1-P
2	C	459	SEP	CA-CB-OG-P
2	C	461	TPO	C-CA-CB-CG2

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Mol	Chain	Res	Type	Atoms
2	C	461	TPO	CB-OG1-P-O2P
2	C	461	TPO	N-CA-CB-CG2

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	461	TPO	3	0
2	C	461	TPO	4	0
2	C	459	SEP	6	0
2	D	459	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	122/122 (100%)	1.17	23 (18%) 3 3	27, 40, 80, 89	0
1	B	118/122 (96%)	2.14	56 (47%) 0 0	32, 55, 84, 103	0
2	C	19/26 (73%)	3.67	17 (89%) 0 0	76, 95, 117, 119	0
2	D	19/26 (73%)	3.14	16 (84%) 0 0	65, 78, 108, 111	0
All	All	278/296 (93%)	1.89	112 (40%) 0 1	27, 52, 95, 119	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	671	SER	6.8
1	B	626	ILE	6.1
2	C	471	ILE	6.0
1	B	717	VAL	6.0
1	B	668	THR	5.9
1	B	718	ALA	5.8
1	A	626	ILE	5.7
2	C	470	PHE	5.6
2	C	477	ALA	5.6
1	B	622	PHE	5.5
1	B	645	THR	5.5
2	C	475	PHE	5.2
2	D	462	VAL	5.1
1	B	646	GLU	5.1
2	D	477	ALA	5.1
2	C	462	VAL	4.9
2	D	457	GLN	4.8
2	C	467	ALA	4.7
1	B	667	ALA	4.6
1	A	622	PHE	4.6
1	B	613	LEU	4.6

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Mol	Chain	Res	Type	RSRZ
2	D	458	PRO	4.6
2	C	458	PRO	4.5
1	A	624	THR	4.4
1	B	674	LYS	4.3
1	B	614	LYS	4.3
1	B	659	PRO	4.2
1	B	643	ILE	4.2
1	B	615	ASN	4.1
1	B	621	VAL	4.0
1	B	716	THR	4.0
2	D	470	PHE	4.0
2	D	460	PRO	3.9
2	D	471	ILE	3.8
1	A	673	SER	3.8
2	D	475	PHE	3.8
2	C	460	PRO	3.8
1	B	618	LEU	3.7
2	C	468	LEU	3.7
1	B	676	GLN	3.7
1	A	618	LEU	3.7
1	B	673	SER	3.7
2	C	474	MET	3.6
1	B	647	ASN	3.6
1	A	621	VAL	3.6
1	B	663	LEU	3.6
1	A	615	ASN	3.6
1	B	649	TYR	3.6
1	B	617	ARG	3.5
1	B	627	GLN	3.5
1	B	619	LYS	3.4
2	C	457	GLN	3.4
1	A	672	GLY	3.3
2	C	472	MET	3.3
1	B	684	HIS	3.2
1	B	650	ARG	3.2
2	D	474	MET	3.2
1	B	665	PHE	3.1
1	B	641	ILE	3.1
1	B	664	ILE	3.1
1	B	640	GLN	3.0
1	A	627	GLN	3.0
1	A	675	MET	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	680	THR	3.0
1	B	610	SER	3.0
1	A	623	GLN	2.9
1	B	612	GLU	2.9
2	D	468	LEU	2.9
1	A	606	LYS	2.8
1	A	631	LYS	2.8
2	D	467	ALA	2.7
1	B	686	VAL	2.7
1	B	644	THR	2.7
1	B	675	MET	2.7
1	A	629	PHE	2.7
2	C	464	THR	2.7
1	A	619	LYS	2.7
2	C	476	GLN	2.6
1	B	651	LEU	2.6
1	A	625	LYS	2.6
1	B	616	GLN	2.5
1	B	611	ALA	2.5
1	B	677	LEU	2.4
2	D	463	HIS	2.4
1	B	648	GLN	2.4
1	B	634	TYR	2.4
1	A	613	LEU	2.4
2	D	472	MET	2.4
1	B	715	GLN	2.3
1	B	694	LEU	2.3
1	B	658	HIS	2.3
2	C	469	GLY	2.3
2	C	463	HIS	2.3
1	B	630	ARG	2.3
1	B	608	VAL	2.3
1	B	624	THR	2.2
1	B	678	LEU	2.2
1	A	597	SER	2.2
1	B	679	GLU	2.2
2	D	476	GLN	2.2
1	B	695	ARG	2.2
2	D	464	THR	2.2
1	B	633	CYS	2.2
1	A	667	ALA	2.1
2	D	473	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	681	GLU	2.1
1	A	611	ALA	2.1
1	A	658	HIS	2.1
1	B	688	GLU	2.1
2	C	473	ASN	2.0
1	A	617	ARG	2.0
1	B	682	PHE	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
2	SEP	C	459	10/11	0.33	0.13	119,125,143,147	0
2	SEP	D	459	10/11	0.50	0.13	107,110,119,121	0
2	TPO	C	461	11/12	0.73	0.14	98,103,105,106	0
2	TPO	D	461	11/12	0.73	0.13	86,93,97,99	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.