



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:09 AM UTC

PDB ID : 9FVL / pdb_00009fvl
Title : Dimeric 14-3-3 zeta in complex with unphosphorylated MAP2c peptide
Authors : Narasimhan, S.; Jansen, S.; Zidek, L.
Deposited on : 2024-06-27
Resolution : 2.08 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
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

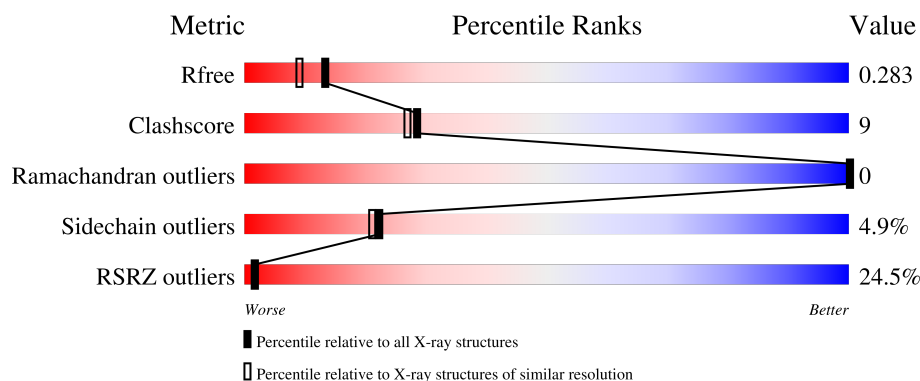
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	248	
1	B	248	
2	E	29	
2	F	29	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3989 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 14-3-3 protein zeta/delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1834	1149	308	368	9			
1	B	225	Total	C	N	O	S	0	0	0
			1797	1127	304	357	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP P63104
A	2	VAL	-	expression tag	UNP P63104
A	3	ASP	-	expression tag	UNP P63104
B	1	SER	-	expression tag	UNP P63104
B	2	VAL	-	expression tag	UNP P63104
B	3	ASP	-	expression tag	UNP P63104

- Molecule 2 is a protein called Isoform 3 of Microtubule-associated protein 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	6	Total	C	N	O	0	0	0
			54	34	13	7			
2	F	6	Total	C	N	O	0	0	0
			54	34	13	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	SER	deletion	UNP P15146
F	?	-	SER	deletion	UNP P15146

- Molecule 3 is CADMIUM ION (CCD ID: CD) (formula: Cd).

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

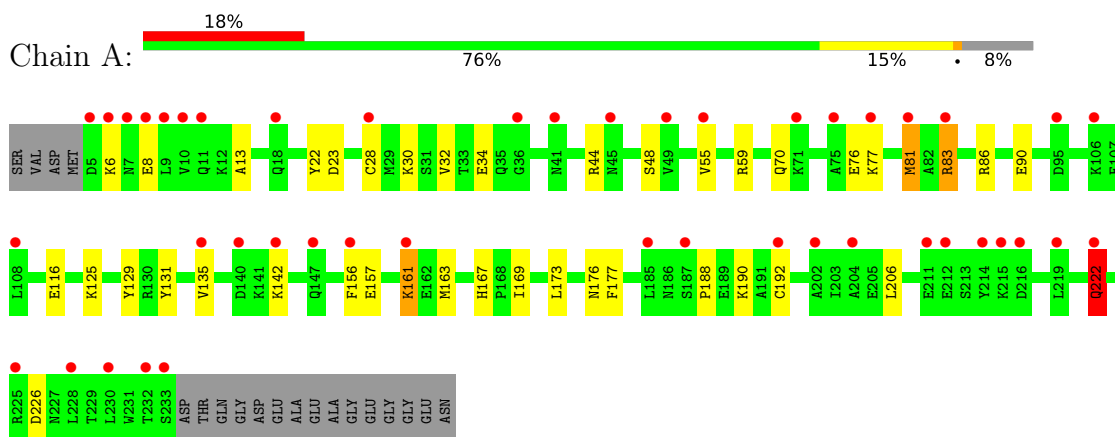
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	125	Total 125	O 125	0	0
4	B	117	Total 117	O 117	0	0
4	E	1	Total 1	O 1	0	0
4	F	2	Total 2	O 2	0	0

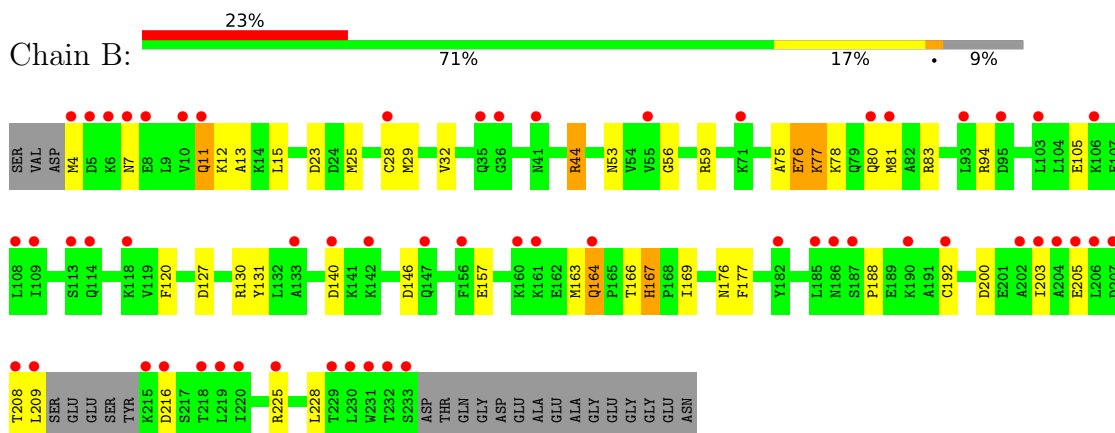
3 Residue-property plots

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.

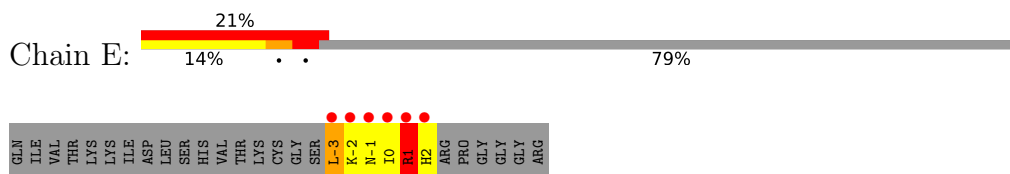
- Molecule 1: 14-3-3 protein zeta/delta



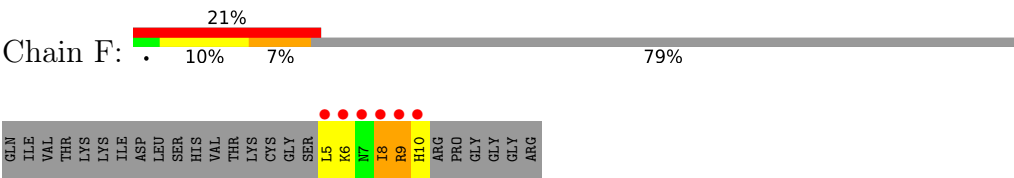
- Molecule 1: 14-3-3 protein zeta/delta



- Molecule 2: Isoform 3 of Microtubule-associated protein 2



- Molecule 2: Isoform 3 of Microtubule-associated protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.31Å 69.11Å 82.62Å 90.00° 112.09° 90.00°	Depositor
Resolution (Å)	48.78 – 2.08 48.78 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.78-2.08) 99.2 (48.78-2.08)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.234 , 0.283 0.240 , 0.283	Depositor DCC
R_{free} test set	2446 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.239	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3989	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.40% 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, 0.333 respectively for untwinned datasets, and 0.375, 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:
CD

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	0.64	0/1859	1.25	6/2500 (0.2%)
1	B	0.64	0/1820	1.29	10/2446 (0.4%)
2	E	0.59	0/54	1.75	2/70 (2.9%)
2	F	0.57	0/54	0.85	0/70
All	All	0.64	0/3787	1.27	18/5086 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
2	E	0	1
2	F	0	1
All	All	0	7

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	76	GLU	CB-CA-C	9.30	128.52	110.67
2	E	1	ARG	CB-CA-C	8.88	122.06	109.71
1	B	11	GLN	CB-CA-C	6.88	122.55	110.85
2	E	1	ARG	N-CA-C	-6.62	100.75	110.42
1	A	76	GLU	CB-CG-CD	6.44	123.55	112.60
1	B	32	VAL	N-CA-CB	6.40	119.25	110.54
1	B	76	GLU	CB-CA-C	6.09	122.81	110.38
1	A	32	VAL	N-CA-CB	5.99	117.95	110.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	PHE	CA-CB-CG	5.89	119.69	113.80
1	B	164	GLN	N-CA-CB	5.83	117.91	110.04
1	B	11	GLN	N-CA-CB	-5.78	101.61	110.16
1	A	222	GLN	CB-CA-C	5.70	119.83	110.88
1	B	167	HIS	CB-CA-C	5.64	118.13	109.50
1	B	23	ASP	CA-CB-CG	5.51	118.11	112.60
1	B	177	PHE	CA-CB-CG	5.51	119.31	113.80
1	B	12	LYS	N-CA-CB	-5.37	102.22	110.01
1	B	164	GLN	CB-CA-C	5.28	117.03	109.11
1	A	23	ASP	CA-CB-CG	5.03	117.63	112.60

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	44	ARG	Sidechain
1	A	83	ARG	Sidechain
1	B	44	ARG	Sidechain
1	B	75	ALA	Peptide
1	B	94	ARG	Sidechain
2	E	1	ARG	Sidechain
2	F	9	ARG	Sidechain

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	1834	0	1821	30	0
1	B	1797	0	1791	31	2
2	E	54	0	63	11	0
2	F	54	0	60	10	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
4	A	125	0	0	9	1
4	B	117	0	0	3	0
4	E	1	0	0	0	0
4	F	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3989	0	3735	66	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 9.

All (66) 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:13:ALA:HB2	1:B:28:CYS:SG	2.12	0.89
1:B:76:GLU:O	1:B:80:GLN:HG2	1.73	0.88
1:B:25:MET:O	1:B:28:CYS:SG	2.33	0.84
1:A:176:ASN:HD21	2:E:2:HIS:HA	1.42	0.84
1:A:59:ARG:HH11	2:E:0:ILE:HG13	1.46	0.80
1:A:90:GLU:HG2	1:A:135:VAL:HG21	1.62	0.80
1:A:8:GLU:OE1	4:A:401:HOH:O	2.03	0.76
1:A:34:GLU:OE2	4:A:402:HOH:O	2.05	0.74
1:B:4:MET:CB	4:B:472:HOH:O	2.36	0.73
1:A:90:GLU:HG2	1:A:135:VAL:CG2	2.20	0.71
1:B:157:GLU:OE2	4:B:401:HOH:O	2.09	0.69
1:B:13:ALA:CB	1:B:28:CYS:SG	2.81	0.69
1:A:188:PRO:O	1:A:192:CYS:SG	2.49	0.68
1:B:131:TYR:OH	2:F:9:ARG:O	2.11	0.68
1:A:90:GLU:CG	1:A:135:VAL:CG2	2.77	0.62
2:E:0:ILE:O	2:E:1:ARG:C	2.43	0.60
1:A:90:GLU:CG	1:A:135:VAL:HG21	2.32	0.60
1:B:225:ARG:HG2	1:B:225:ARG:HH11	1.67	0.59
1:A:59:ARG:NH1	2:E:0:ILE:HG13	2.16	0.59
1:A:90:GLU:CD	1:A:135:VAL:CG2	2.75	0.59
1:B:59:ARG:HD3	2:F:8:ILE:HG21	1.85	0.59
1:B:80:GLN:NE2	1:B:83:ARG:HH21	2.01	0.59
4:A:407:HOH:O	2:F:6:LYS:NZ	2.36	0.58
1:B:80:GLN:NE2	1:B:83:ARG:NH2	2.52	0.57
1:A:59:ARG:HH11	2:E:0:ILE:CG1	2.14	0.56
1:A:176:ASN:HD21	2:E:2:HIS:CA	2.17	0.56
1:A:125:LYS:HE2	1:A:129:TYR:HE2	1.71	0.55
1:B:188:PRO:O	1:B:192:CYS:SG	2.59	0.53
1:A:157:GLU:OE2	4:A:403:HOH:O	2.19	0.52
1:A:222:GLN:NE2	1:A:226:ASP:OD2	2.43	0.51
1:A:131:TYR:OH	2:E:1:ARG:O	2.29	0.51
1:A:30:LYS:NZ	4:A:402:HOH:O	2.42	0.51
1:B:28:CYS:SG	1:B:29:MET:N	2.84	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:ASP:O	1:B:203:ILE:HG22	2.10	0.51
1:A:83:ARG:NH1	4:A:410:HOH:O	2.44	0.50
1:A:156:PHE:HA	1:A:173:LEU:HD21	1.94	0.48
1:B:208:THR:O	1:B:209:LEU:C	2.56	0.48
1:A:81:MET:HE1	4:A:493:HOH:O	2.14	0.47
1:B:77:LYS:HE2	1:B:78:LYS:H	1.80	0.46
1:A:116:GLU:HA	1:A:163:MET:HE3	1.97	0.46
1:B:56:GLY:HA2	2:F:8:ILE:CG2	2.46	0.46
1:B:53:ASN:HA	2:F:6:LYS:O	2.16	0.45
2:E:-3:LEU:HD13	2:F:5:LEU:CD1	2.46	0.45
1:B:80:GLN:HE22	1:B:83:ARG:NH2	2.13	0.45
1:B:164:GLN:OE1	1:B:205:GLU:OE2	2.33	0.45
1:B:77:LYS:HD2	1:B:78:LYS:N	2.31	0.45
1:B:127:ASP:OD2	2:F:10:HIS:ND1	2.50	0.44
1:A:163:MET:HE2	1:A:167:HIS:CE1	2.53	0.43
2:E:1:ARG:O	2:E:2:HIS:ND1	2.52	0.43
1:A:22:TYR:HD2	2:E:-3:LEU:HD23	1.84	0.43
1:B:44:ARG:HD2	1:B:120:PHE:CD1	2.54	0.43
1:A:161:LYS:O	1:A:161:LYS:HD3	2.19	0.42
1:B:7:ASN:HB3	4:B:459:HOH:O	2.18	0.42
2:E:0:ILE:HD12	2:E:0:ILE:HA	1.97	0.42
1:A:142:LYS:HB3	4:A:469:HOH:O	2.20	0.41
1:A:167:HIS:HE1	1:A:169:ILE:HD12	1.84	0.41
1:A:86:ARG:O	1:A:90:GLU:HG3	2.20	0.41
1:B:176:ASN:HD21	2:F:10:HIS:HA	1.85	0.41
1:B:225:ARG:HG2	1:B:225:ARG:NH1	2.34	0.41
1:A:13:ALA:HB2	1:A:28:CYS:SG	2.60	0.41
1:A:90:GLU:CD	1:A:135:VAL:HG22	2.46	0.41
4:A:459:HOH:O	1:B:81:MET:HE1	2.21	0.41
1:B:59:ARG:HH11	2:F:8:ILE:CG2	2.35	0.40
1:B:163:MET:HE1	1:B:169:ILE:HB	2.02	0.40
1:B:130:ARG:NH2	2:F:9:ARG:O	2.52	0.40
1:B:166:THR:O	1:B:167:HIS:C	2.65	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:76:GLU:OE1	1:B:105:GLU:OE2[2_345]	2.07	0.13
1:B:146:ASP:OD2	4:A:401:HOH:O[2_355]	2.14	0.06

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	227/248 (92%)	225 (99%)	2 (1%)	0	100	100
1	B	221/248 (89%)	216 (98%)	5 (2%)	0	100	100
2	E	4/29 (14%)	2 (50%)	2 (50%)	0	100	100
2	F	4/29 (14%)	2 (50%)	2 (50%)	0	100	100
All	All	456/554 (82%)	445 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	199/212 (94%)	189 (95%)	10 (5%)	22	20
1	B	194/212 (92%)	188 (97%)	6 (3%)	35	38
2	E	6/25 (24%)	3 (50%)	3 (50%)	0	0
2	F	6/25 (24%)	5 (83%)	1 (17%)	2	1
All	All	405/474 (85%)	385 (95%)	20 (5%)	22	21

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	48	SER

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Mol	Chain	Res	Type
1	A	55	VAL
1	A	70	GLN
1	A	77	LYS
1	A	81	MET
1	A	161	LYS
1	A	190	LYS
1	A	206	LEU
1	A	222	GLN
1	B	11	GLN
1	B	15	LEU
1	B	77	LYS
1	B	140	ASP
1	B	216	ASP
1	B	228	LEU
2	E	-3	LEU
2	E	-2	LYS
2	E	-1	ASN
2	F	8	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	176	ASN
1	A	186	ASN
1	B	53	ASN
1	B	80	GLN
1	B	150	GLN
2	F	7	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 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	229/248 (92%)	1.19	45 (19%) 3 3	22, 34, 59, 89	0
1	B	225/248 (90%)	1.34	57 (25%) 1 1	23, 35, 64, 89	0
2	E	6/29 (20%)	6.91	6 (100%) 0 0	69, 84, 89, 90	0
2	F	6/29 (20%)	7.00	6 (100%) 0 0	72, 80, 85, 87	0
All	All	466/554 (84%)	1.41	114 (24%) 2 2	22, 35, 69, 90	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	-3	LEU	10.7
2	F	5	LEU	10.1
2	F	6	LYS	8.0
2	E	-2	LYS	7.7
1	A	192	CYS	7.6
2	F	10	HIS	7.4
2	F	8	ILE	7.2
2	E	2	HIS	7.1
2	E	0	ILE	7.0
1	B	215	LYS	6.8
1	B	4	MET	6.7
1	B	192	CYS	5.9
1	A	28	CYS	5.7
1	B	209	LEU	5.6
2	F	7	ASN	5.2
1	B	232	THR	5.0
1	B	204	ALA	4.8
2	E	1	ARG	4.6
2	E	-1	ASN	4.5
1	B	5	ASP	4.2
2	F	9	ARG	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	28	CYS	4.0
1	B	207	ASP	4.0
1	B	161	LYS	3.9
1	A	5	ASP	3.9
1	B	218	THR	3.7
1	A	215	LYS	3.7
1	A	216	ASP	3.6
1	B	156	PHE	3.5
1	B	216	ASP	3.5
1	A	232	THR	3.5
1	A	230	LEU	3.4
1	B	203	ILE	3.4
1	A	142	LYS	3.3
1	B	142	LYS	3.3
1	A	49	VAL	3.2
1	B	160	LYS	3.2
1	A	8	GLU	3.2
1	B	80	GLN	3.2
1	B	208	THR	3.1
1	A	81	MET	3.1
1	B	230	LEU	3.1
1	B	202	ALA	3.1
1	A	6	LYS	3.1
1	A	233	SER	3.1
1	B	147	GLN	3.0
1	A	108	LEU	3.0
1	B	164	GLN	2.9
1	B	11	GLN	2.8
1	B	103	LEU	2.8
1	A	7	ASN	2.8
1	B	93	LEU	2.7
1	A	10	VAL	2.7
1	B	10	VAL	2.7
1	B	81	MET	2.7
1	A	11	GLN	2.7
1	A	95	ASP	2.7
1	A	106	LYS	2.7
1	A	9	LEU	2.7
1	B	206	LEU	2.7
1	A	161	LYS	2.6
1	A	41	ASN	2.6
1	A	147	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	35	GLN	2.5
1	B	41	ASN	2.5
1	A	55	VAL	2.5
1	B	71	LYS	2.5
1	B	231	TRP	2.5
1	B	36	GLY	2.5
1	A	83	ARG	2.5
1	B	182	TYR	2.5
1	A	135	VAL	2.4
1	A	222	GLN	2.4
1	B	233	SER	2.4
1	A	204	ALA	2.4
1	B	185	LEU	2.4
1	A	225	ARG	2.4
1	A	156	PHE	2.4
1	B	140	ASP	2.4
1	A	187	SER	2.4
1	B	219	LEU	2.4
1	B	8	GLU	2.4
1	B	109	ILE	2.4
1	B	6	LYS	2.4
1	A	219	LEU	2.4
1	B	190	LYS	2.4
1	B	133	ALA	2.3
1	A	228	LEU	2.3
1	A	36	GLY	2.3
1	B	106	LYS	2.3
1	B	113	SER	2.3
1	B	229	THR	2.3
1	B	187	SER	2.3
1	A	140	ASP	2.3
1	A	211	GLU	2.2
1	A	212	GLU	2.2
1	B	55	VAL	2.2
1	B	205	GLU	2.2
1	A	18	GLN	2.2
1	B	7	ASN	2.2
1	B	186	ASN	2.2
1	A	202	ALA	2.1
1	A	185	LEU	2.1
1	B	118	LYS	2.1
1	B	225	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	75	ALA	2.1
1	B	114	GLN	2.1
1	B	108	LEU	2.1
1	A	71	LYS	2.1
1	B	220	ILE	2.1
1	A	45	ASN	2.0
1	A	77	LYS	2.0
1	A	214	TYR	2.0
1	B	95	ASP	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 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($\text{\AA}^2$)	Q<0.9
3	CD	A	303	1/1	0.77	0.12	96,96,96,96	0
3	CD	A	302	1/1	0.99	0.02	33,33,33,33	0
3	CD	A	301	1/1	0.99	0.04	32,32,32,32	0
3	CD	B	302	1/1	0.99	0.05	33,33,33,33	0
3	CD	B	301	1/1	1.00	0.03	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.