



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2017 – 04:16 AM EDT

PDB ID : 3KMD
Title : Crystal structure of the p53 core domain bound to a full consensus site as a self-assembled tetramer
Authors : Chen, Y.; Dey, R.; Chen, L.
Deposited on : unknown
Resolution : 2.15 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030345

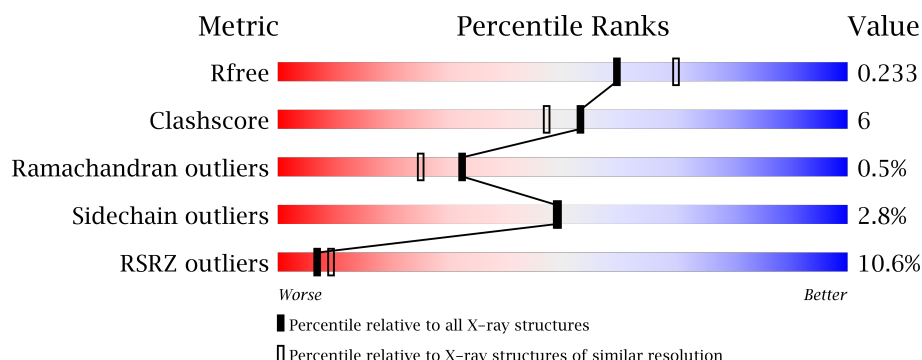
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	100719	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

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	200	<div> <div>7%</div> <div>90%</div> <div>9%</div> <div>.</div> </div>
1	B	200	<div> <div>10%</div> <div>85%</div> <div>14%</div> <div>.</div> </div>
1	C	200	<div> <div>19%</div> <div>87%</div> <div>12%</div> <div>..</div> </div>
1	D	200	<div> <div>8%</div> <div>89%</div> <div>12%</div> </div>
2	E	19	<div> <div>5%</div> <div>53%</div> <div>37%</div> <div>11%</div> </div>

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Mol	Chain	Length	Quality of chain	
				
2	F	19	53%	37%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7587 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 Cellular tumor antigen p53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1571	968	291	296	16			
1	B	200	Total	C	N	O	S	0	0	0
			1571	968	291	296	16			
1	D	200	Total	C	N	O	S	0	0	0
			1571	968	291	296	16			
1	C	200	Total	C	N	O	S	0	0	0
			1571	968	291	296	16			

- Molecule 2 is a DNA chain called 5'-D(*GP*GP*GP*CP*AP*TP*GP*CP*CP*TP*AP*GP*GP*CP*AP*TP*GP*CP*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	19	Total	C	N	O	P	0	0	0
			388	184	74	112	18			
2	F	19	Total	C	N	O	P	0	0	0
			388	184	74	112	18			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

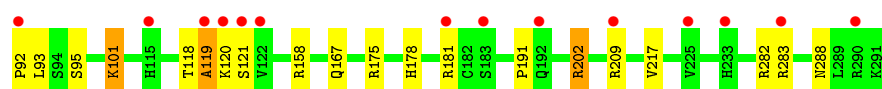
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	156	Total 156	O 156	0	0
4	B	84	Total 84	O 84	0	0
4	D	128	Total 128	O 128	0	0
4	C	101	Total 101	O 101	0	0
4	E	30	Total 30	O 30	0	0
4	F	24	Total 24	O 24	0	0

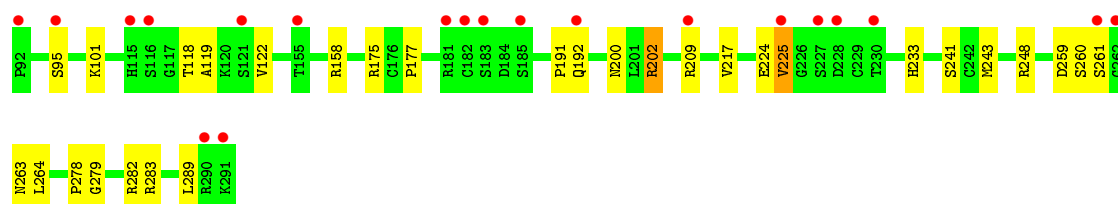
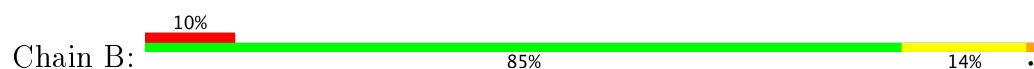
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

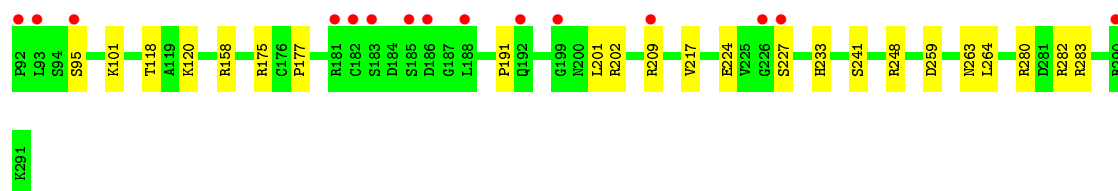
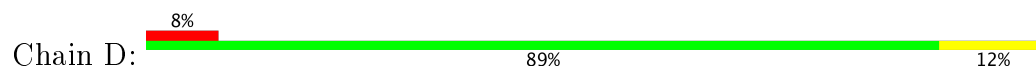
• Molecule 1: Cellular tumor antigen p53



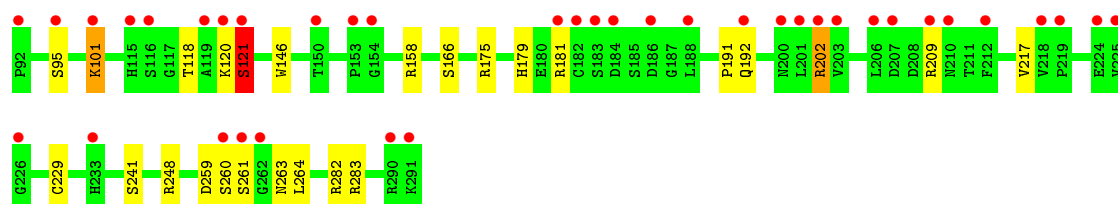
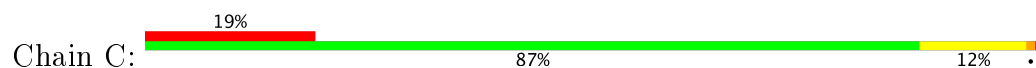
• Molecule 1: Cellular tumor antigen p53



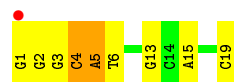
• Molecule 1: Cellular tumor antigen p53



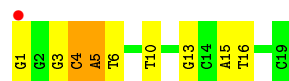
• Molecule 1: Cellular tumor antigen p53



- Molecule 2: 5'-D(*GP*GP*GP*CP*AP*TP*GP*CP*CP*TP*AP*GP*GP*CP*AP*TP*GP*CP*C)-3'



- Molecule 2: 5'-D(*GP*GP*GP*CP*AP*TP*GP*CP*CP*TP*AP*GP*GP*CP*AP*TP*GP*CP*C)-3'



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.08 Å 93.71 Å 145.57 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.52 – 2.15 48.52 – 2.06	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.52-2.15) 99.2 (48.52-2.06)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.07 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.216 , 0.239 0.212 , 0.233	Depositor DCC
R_{free} test set	2524 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.441	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7587	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.87 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.5674e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 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	0.39	0/1607	0.56	1/2176 (0.0%)
1	B	0.34	0/1607	0.51	0/2176
1	C	0.34	0/1607	0.53	1/2176 (0.0%)
1	D	0.38	0/1607	0.55	1/2176 (0.0%)
2	E	0.67	0/435	1.50	6/670 (0.9%)
2	F	0.76	0/435	1.56	10/670 (1.5%)
All	All	0.42	0/7298	0.75	19/10044 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	5	DA	O4'-C1'-N9	-10.70	100.51	108.00
2	E	13	DG	O4'-C1'-N9	9.76	114.83	108.00
2	F	15	DA	O4'-C1'-N9	-9.71	101.21	108.00
2	E	15	DA	O4'-C1'-N9	-9.46	101.38	108.00
2	F	5	DA	O4'-C1'-N9	-7.72	102.60	108.00
2	F	6	DT	C1'-O4'-C4'	-6.86	103.25	110.10
2	E	13	DG	C3'-C2'-C1'	-6.68	94.49	102.50
2	F	16	DT	O4'-C1'-C2'	-5.97	101.13	105.90
1	A	119	ALA	N-CA-C	-5.95	94.93	111.00
2	E	4	DC	O4'-C1'-N1	5.91	112.14	108.00
1	D	227	SER	CB-CA-C	-5.81	99.06	110.10
2	F	4	DC	C4'-C3'-C2'	-5.79	97.89	103.10
2	F	16	DT	P-O3'-C3'	5.77	126.62	119.70
1	C	121	SER	CB-CA-C	5.72	120.96	110.10
2	F	1	DG	P-O3'-C3'	5.49	126.29	119.70
2	E	6	DT	C1'-O4'-C4'	-5.32	104.78	110.10
2	F	16	DT	C1'-O4'-C4'	-5.24	104.86	110.10
2	F	13	DG	C3'-C2'-C1'	-5.22	96.24	102.50

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	F	13	DG	O4'-C1'-N9	5.13	111.59	108.00

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	1571	0	1535	21	0
1	B	1571	0	1535	27	0
1	C	1571	0	1535	27	0
1	D	1571	0	1535	17	0
2	E	388	0	214	5	0
2	F	388	0	214	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	156	0	0	3	0
4	B	84	0	0	1	0
4	C	101	0	0	2	0
4	D	128	0	0	2	0
4	E	30	0	0	1	0
4	F	24	0	0	2	0
All	All	7587	0	6568	85	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 6.

All (85) 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:202:ARG:HH11	1:A:202:ARG:HG3	1.19	1.05
1:C:202:ARG:HH11	1:C:202:ARG:HG3	1.20	1.04
1:B:202:ARG:HH11	1:B:202:ARG:HG3	1.20	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:GLU:HG3	1:C:101:LYS:HD3	1.43	1.01
2:F:3:DG:H2"	2:F:4:DC:H5"	1.53	0.91
1:B:202:ARG:HG3	1:B:202:ARG:NH1	1.98	0.78
1:B:224:GLU:CG	1:C:101:LYS:HD3	2.16	0.75
1:D:202:ARG:HH11	1:D:202:ARG:HG3	1.56	0.71
1:A:101:LYS:HD3	1:D:224:GLU:HB3	1.71	0.71
2:E:2:DG:H1'	2:E:3:DG:H5"	1.73	0.69
1:A:202:ARG:NH1	1:A:202:ARG:HG3	1.98	0.68
1:B:202:ARG:HH11	1:B:202:ARG:CG	2.04	0.68
1:B:118:THR:HG22	1:B:282:ARG:HD3	1.77	0.65
1:A:202:ARG:CG	1:A:202:ARG:HH11	2.03	0.65
1:A:118:THR:OG1	1:A:283:ARG:HD3	1.97	0.65
1:D:202:ARG:NH1	4:D:447:HOH:O	2.30	0.63
1:A:167:GLN:HB3	1:D:233:HIS:CD2	2.33	0.62
1:A:181:ARG:HD2	1:B:177:PRO:HB3	1.81	0.62
1:C:120:LYS:HB3	2:E:1:DG:H8	1.64	0.62
1:C:118:THR:HG22	1:C:282:ARG:HD3	1.82	0.62
1:D:118:THR:HG22	1:D:282:ARG:HD3	1.82	0.61
1:C:202:ARG:NH1	1:C:202:ARG:HG3	1.99	0.61
1:A:119:ALA:O	1:A:121:SER:N	2.33	0.60
1:A:101:LYS:HB2	1:D:224:GLU:OE2	2.01	0.60
1:B:241:SER:HA	1:B:248:ARG:HG3	1.85	0.58
1:A:181:ARG:HH11	1:B:177:PRO:HG3	1.67	0.58
2:F:3:DG:C2'	2:F:4:DC:H5"	2.33	0.56
1:C:118:THR:OG1	1:C:283:ARG:HD3	2.06	0.55
2:F:5:DA:H5'	4:F:479:HOH:O	2.06	0.55
1:C:120:LYS:O	1:C:121:SER:HB2	2.06	0.55
1:C:175:ARG:HD3	1:C:191:PRO:O	2.07	0.55
1:D:118:THR:OG1	1:D:283:ARG:HD3	2.07	0.54
1:C:202:ARG:HH11	1:C:202:ARG:CG	2.04	0.54
2:F:4:DC:H2"	2:F:5:DA:C8	2.43	0.54
1:D:202:ARG:NH1	1:D:202:ARG:HG3	2.19	0.53
1:A:92:PRO:N	4:A:450:HOH:O	2.41	0.53
2:E:19:DC:H2"	4:E:373:HOH:O	2.08	0.53
1:D:241:SER:HA	1:D:248:ARG:HG3	1.90	0.52
1:C:241:SER:HA	1:C:248:ARG:HG3	1.89	0.52
1:C:259:ASP:OD1	1:C:263:ASN:HB2	2.10	0.52
1:B:175:ARG:HD3	1:B:191:PRO:O	2.10	0.51
1:B:119:ALA:O	1:B:279:GLY:HA3	2.10	0.51
1:B:202:ARG:NH1	1:B:202:ARG:CG	2.69	0.51
1:B:225:VAL:HB	1:C:101:LYS:HE2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:ASP:OD2	1:B:263:ASN:HB2	2.12	0.50
1:A:175:ARG:HD3	1:A:191:PRO:O	2.12	0.50
1:A:158:ARG:HG3	1:A:217:VAL:HG22	1.93	0.49
1:C:202:ARG:NH1	1:C:202:ARG:CG	2.69	0.49
1:D:177:PRO:HG3	1:C:181:ARG:HH11	1.77	0.48
2:E:4:DC:H2''	2:E:5:DA:C8	2.48	0.48
1:D:158:ARG:CZ	1:D:217:VAL:HG21	2.43	0.48
1:B:233:HIS:CE1	1:C:166:SER:OG	2.66	0.48
1:C:120:LYS:HB3	2:E:1:DG:C8	2.48	0.47
1:D:175:ARG:HD3	1:D:191:PRO:O	2.13	0.47
1:A:118:THR:HG22	1:A:282:ARG:HD3	1.97	0.47
1:A:202:ARG:CG	1:A:202:ARG:NH1	2.68	0.47
1:B:118:THR:OG1	1:B:283:ARG:HD3	2.15	0.46
1:A:158:ARG:CZ	1:A:217:VAL:HG21	2.45	0.46
1:B:225:VAL:HB	1:C:101:LYS:CE	2.46	0.46
1:C:259:ASP:O	1:C:261:SER:N	2.49	0.46
2:F:10:DT:H5'	4:F:471:HOH:O	2.16	0.46
1:A:92:PRO:CA	4:A:450:HOH:O	2.63	0.46
1:B:200:ASN:HA	4:B:322:HOH:O	2.15	0.45
1:C:120:LYS:O	1:C:121:SER:CB	2.65	0.44
1:D:158:ARG:HG3	1:D:217:VAL:HG22	1.98	0.44
1:A:178:HIS:CD2	1:B:243:MET:HG3	2.53	0.44
1:C:158:ARG:CZ	1:C:217:VAL:HG21	2.48	0.43
1:B:224:GLU:O	1:B:225:VAL:C	2.57	0.43
1:D:120:LYS:HD2	1:D:280:ARG:HB2	2.02	0.42
1:B:158:ARG:CZ	1:B:217:VAL:HG21	2.49	0.42
1:C:146:TRP:CE2	1:C:229:CYS:HB3	2.54	0.42
1:C:248:ARG:NH2	4:C:350:HOH:O	2.53	0.42
1:B:192:GLN:H	1:B:192:GLN:HG2	1.70	0.42
1:B:289:LEU:HD12	1:B:289:LEU:HA	1.93	0.41
1:C:192:GLN:NE2	4:C:458:HOH:O	2.52	0.41
1:A:93:LEU:HB2	1:D:201:LEU:HG	2.02	0.41
1:C:175:ARG:NH2	1:C:179:HIS:HB3	2.35	0.41
1:D:259:ASP:OD2	1:D:263:ASN:HB2	2.20	0.41
1:A:178:HIS:CG	1:B:243:MET:HG3	2.54	0.41
1:B:259:ASP:CG	1:B:263:ASN:HB2	2.41	0.41
1:A:288:ASN:ND2	4:A:330:HOH:O	2.54	0.41
1:B:122:VAL:O	1:B:278:PRO:HG2	2.20	0.40
1:C:158:ARG:HG3	1:C:217:VAL:HG22	2.03	0.40
1:D:118:THR:HG21	4:D:305:HOH:O	2.21	0.40
1:B:225:VAL:CG2	1:C:101:LYS:HE2	2.51	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	198/200 (99%)	194 (98%)	3 (2%)	1 (0%)	32	25
1	B	198/200 (99%)	193 (98%)	4 (2%)	1 (0%)	32	25
1	C	198/200 (99%)	193 (98%)	3 (2%)	2 (1%)	18	11
1	D	198/200 (99%)	193 (98%)	5 (2%)	0	100	100
All	All	792/800 (99%)	773 (98%)	15 (2%)	4 (0%)	32	25

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	LYS
1	C	260	SER
1	C	121	SER
1	B	225	VAL

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	180/180 (100%)	176 (98%)	4 (2%)	57	60
1	B	180/180 (100%)	173 (96%)	7 (4%)	37	34
1	C	180/180 (100%)	175 (97%)	5 (3%)	49	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	180/180 (100%)	176 (98%)	4 (2%)	57	60
All	All	720/720 (100%)	700 (97%)	20 (3%)	49	49

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

Mol	Chain	Res	Type
1	A	95	SER
1	A	101	LYS
1	A	202	ARG
1	A	209	ARG
1	B	95	SER
1	B	101	LYS
1	B	202	ARG
1	B	209	ARG
1	B	260	SER
1	B	261	SER
1	B	264	LEU
1	D	95	SER
1	D	101	LYS
1	D	209	ARG
1	D	264	LEU
1	C	95	SER
1	C	101	LYS
1	C	202	ARG
1	C	209	ARG
1	C	264	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	247	ASN
1	A	263	ASN
1	A	288	ASN
1	B	100	GLN
1	B	247	ASN
1	B	288	ASN
1	D	247	ASN
1	D	288	ASN
1	C	100	GLN
1	C	178	HIS
1	C	247	ASN

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Mol	Chain	Res	Type
1	C	263	ASN
1	C	268	ASN
1	C	288	ASN

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 [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	200/200 (100%)	0.41	14 (7%) 17 22	12, 24, 52, 115	0
1	B	200/200 (100%)	0.72	20 (10%) 8 11	16, 33, 83, 105	0
1	C	200/200 (100%)	0.92	38 (19%) 1 2	17, 35, 85, 140	0
1	D	200/200 (100%)	0.49	15 (7%) 15 19	13, 25, 72, 107	0
2	E	19/19 (100%)	0.19	1 (5%) 27 34	22, 35, 67, 76	0
2	F	19/19 (100%)	0.19	1 (5%) 27 34	21, 35, 62, 73	0
All	All	838/838 (100%)	0.61	89 (10%) 7 9	12, 29, 74, 140	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	121	SER	7.6
1	D	183	SER	7.1
1	A	115	HIS	6.5
1	A	121	SER	6.2
1	D	92	PRO	6.1
1	D	209	ARG	5.9
1	C	209	ARG	5.9
1	A	120	LYS	5.3
1	B	95	SER	5.1
1	B	92	PRO	4.9
1	A	209	ARG	4.8
1	B	225	VAL	4.7
1	B	209	ARG	4.5
1	C	153	PRO	4.4
1	C	201	LEU	4.3
1	D	181	ARG	4.3
1	B	290	ARG	4.1
1	B	291	LYS	4.0
1	C	181	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	183	SER	3.9
1	C	261	SER	3.9
1	A	181	ARG	3.8
1	D	188	LEU	3.8
1	C	120	LYS	3.7
1	C	210	ASN	3.7
1	C	226	GLY	3.6
1	C	150	THR	3.6
1	C	92	PRO	3.5
1	C	291	LYS	3.5
1	A	290	ARG	3.5
1	B	181	ARG	3.3
1	D	93	LEU	3.2
1	D	290	ARG	3.2
1	C	225	VAL	3.2
1	C	154	GLY	3.1
1	B	230	THR	3.1
1	B	115	HIS	3.0
1	C	183	SER	3.0
1	C	218	VAL	3.0
1	C	115	HIS	3.0
1	C	262	GLY	3.0
1	C	119	ALA	3.0
1	B	228	ASP	2.9
1	C	203	VAL	2.9
1	C	200	ASN	2.9
2	F	1	DG	2.9
1	C	202	ARG	2.9
2	E	1	DG	2.9
1	A	119	ALA	2.8
1	B	227	SER	2.8
1	B	262	GLY	2.8
1	C	184	ASP	2.8
1	D	95	SER	2.8
1	D	185	SER	2.8
1	C	192	GLN	2.7
1	B	185	SER	2.7
1	D	227	SER	2.7
1	D	182	CYS	2.7
1	C	219	PRO	2.6
1	A	225	VAL	2.6
1	D	226	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	192	GLN	2.6
1	C	233	HIS	2.5
1	A	283	ARG	2.5
1	A	183	SER	2.5
1	C	206	LEU	2.5
1	B	192	GLN	2.4
1	C	260	SER	2.4
1	B	155	THR	2.4
1	D	186	ASP	2.4
1	B	182	CYS	2.4
1	A	92	PRO	2.4
1	C	101	LYS	2.3
1	A	122	VAL	2.2
1	C	182	CYS	2.2
1	C	116	SER	2.2
1	C	186	ASP	2.2
1	B	116	SER	2.2
1	B	261	SER	2.2
1	C	95	SER	2.2
1	C	207	ASP	2.1
1	D	199	GLY	2.1
1	C	188	LEU	2.1
1	C	224	GLU	2.1
1	D	192	GLN	2.1
1	C	212	PHE	2.1
1	A	233	HIS	2.0
1	B	121	SER	2.0
1	C	290	ARG	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(Å ²)	Q<0.9
3	ZN	A	1	1/1	0.99	0.14	0.30	21,21,21,21	0
3	ZN	C	1	1/1	1.00	0.10	-0.81	27,27,27,27	0
3	ZN	D	1	1/1	1.00	0.10	-1.30	21,21,21,21	0
3	ZN	B	1	1/1	1.00	0.09	-1.76	19,19,19,19	0

6.5 Other polymers

There are no such residues in this entry.