



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:40 AM GMT

PDB ID : 2NS2
Title : Crystal Structure of Spindlin1
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Deposited on : 2006-11-02
Resolution : 2.20 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

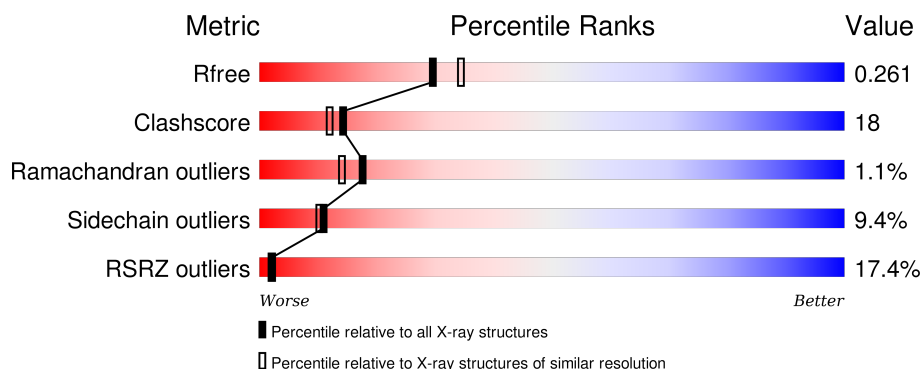
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	242	<div> <div>5%</div> <div>63%</div> <div>17%</div> <div>•</div> <div>17%</div> </div>
1	B	242	<div> <div>21%</div> <div>42%</div> <div>25%</div> <div>•</div> <div>29%</div> </div>

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	PO4	A	301	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3274 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 Spindlin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1626	1039	270	308	9			
1	B	172	Total	C	N	O	S	0	0	0
			1397	903	225	260	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	CLONING ARTIFACT	UNP Q9Y657
A	-3	PRO	-	CLONING ARTIFACT	UNP Q9Y657
A	-2	LEU	-	CLONING ARTIFACT	UNP Q9Y657
A	-1	GLY	-	CLONING ARTIFACT	UNP Q9Y657
A	0	SER	-	CLONING ARTIFACT	UNP Q9Y657
B	-4	GLY	-	CLONING ARTIFACT	UNP Q9Y657
B	-3	PRO	-	CLONING ARTIFACT	UNP Q9Y657
B	-2	LEU	-	CLONING ARTIFACT	UNP Q9Y657
B	-1	GLY	-	CLONING ARTIFACT	UNP Q9Y657
B	0	SER	-	CLONING ARTIFACT	UNP Q9Y657

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

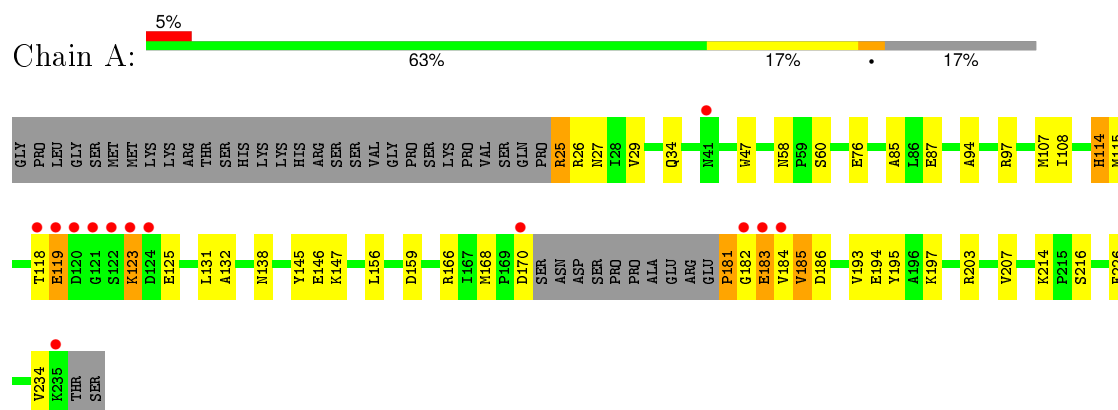
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	170	Total	O	0	0
			170	170		
3	B	71	Total	O	0	0
			71	71		

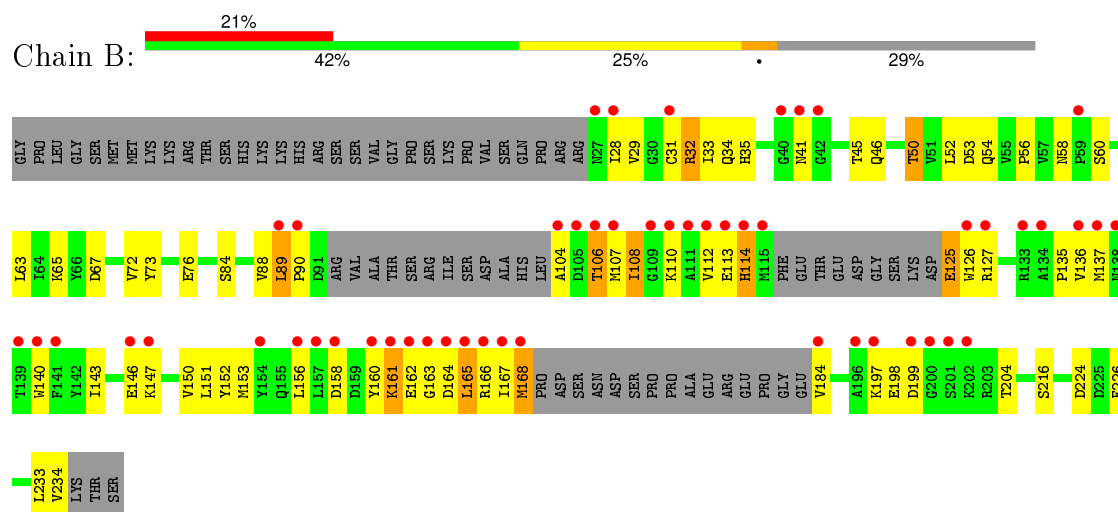
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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: Spindlin-1



• Molecule 1: Spindlin-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	40.79 Å 84.88 Å 136.57 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.38 – 2.20 40.12 – 2.13	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.38-2.20) 98.6 (40.12-2.13)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.75 (at 2.12 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.225 , 0.260 0.226 , 0.261	Depositor DCC
R_{free} test set	2453 reflections (9.97%)	DCC
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 69.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 27366 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3274	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.45	0/1664	0.79	1/2252 (0.0%)
1	B	0.37	0/1429	0.75	0/1934
All	All	0.41	0/3093	0.77	1/4186 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	VAL	N-CA-C	5.25	125.18	111.00

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	1626	0	1585	35	0
1	B	1397	0	1364	76	0
2	A	10	0	0	0	0
3	A	170	0	0	1	0
3	B	71	0	0	0	0
All	All	3274	0	2949	108	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 18.

All (108) 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:185:VAL:HG12	1:A:186:ASP:H	1.22	1.03
1:A:25:ARG:HH11	1:A:25:ARG:HG2	1.28	0.97
1:B:114:HIS:HB2	1:B:164:ASP:O	1.68	0.93
1:B:197:LYS:HG2	1:B:198:GLU:H	1.33	0.92
1:B:65:LYS:HB2	1:B:153:MET:HE3	1.60	0.84
1:B:65:LYS:HD3	1:B:153:MET:HE3	1.59	0.83
1:A:185:VAL:HG12	1:A:186:ASP:N	1.99	0.77
1:B:32:ARG:HE	1:B:50:THR:HG22	1.48	0.77
1:A:125:GLU:OE1	1:A:166:ARG:NH1	2.18	0.76
1:B:127:ARG:HD3	1:B:146:GLU:HG3	1.68	0.75
1:B:158:ASP:O	1:B:161:LYS:HE2	1.89	0.73
1:B:65:LYS:HD3	1:B:153:MET:CE	2.20	0.72
1:B:72:VAL:HG22	1:B:153:MET:HE1	1.70	0.71
1:A:185:VAL:CG1	1:A:186:ASP:H	2.03	0.70
1:B:28:ILE:HG22	1:B:31:CYS:HB2	1.71	0.70
1:B:89:LEU:CG	1:B:90:PRO:HD2	2.22	0.69
1:B:161:LYS:HD2	1:B:161:LYS:C	2.12	0.69
1:B:161:LYS:HD2	1:B:162:GLU:N	2.08	0.69
1:A:25:ARG:NH1	1:A:25:ARG:HG2	2.00	0.68
1:B:127:ARG:CD	1:B:146:GLU:HG3	2.25	0.67
1:B:197:LYS:CG	1:B:198:GLU:H	2.08	0.66
1:B:89:LEU:HG	1:B:90:PRO:HD2	1.77	0.66
1:B:216:SER:OG	1:B:234:VAL:HB	1.97	0.63
1:B:137:MET:HB3	1:B:140:TRP:CD1	2.34	0.62
1:A:182:GLY:O	1:A:183:GLU:HB2	2.00	0.62
1:B:34:GLN:HE21	1:B:46:GLN:CG	2.13	0.61
1:B:89:LEU:CB	1:B:90:PRO:HD2	2.32	0.60
1:B:32:ARG:NH2	1:B:67:ASP:OD1	2.35	0.60
1:B:110:LYS:HD3	1:B:167:ILE:CG2	2.32	0.59
1:B:72:VAL:HG22	1:B:153:MET:CE	2.31	0.59
1:B:65:LYS:CD	1:B:153:MET:HE3	2.31	0.59
1:B:33:ILE:HD12	1:B:35:HIS:HD2	1.68	0.58
1:B:114:HIS:CB	1:B:165:LEU:HD12	2.34	0.58
1:B:137:MET:HB3	1:B:140:TRP:HD1	1.68	0.58
1:B:60:SER:O	1:B:76:GLU:HA	2.04	0.58
1:B:104:ALA:O	1:B:108:ILE:HG13	2.04	0.58
1:A:107:MET:HE1	1:A:156:LEU:HB3	1.86	0.57
1:B:65:LYS:CB	1:B:153:MET:HE3	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:SER:HB2	1:A:234:VAL:HB	1.88	0.55
1:A:146:GLU:HG3	1:A:147:LYS:HG3	1.87	0.55
1:B:161:LYS:C	1:B:163:GLY:H	2.09	0.55
1:B:28:ILE:CG2	1:B:31:CYS:HB2	2.37	0.55
1:B:160:TYR:N	1:B:165:LEU:HD22	2.23	0.54
1:A:114:HIS:HD2	1:A:145:TYR:OH	1.91	0.53
1:A:193:VAL:HG12	1:B:45:THR:HG22	1.91	0.53
1:B:126:TRP:HZ3	1:B:147:LYS:HB2	1.74	0.52
1:A:60:SER:O	1:A:76:GLU:HA	2.10	0.52
1:B:197:LYS:HG2	1:B:198:GLU:N	2.12	0.52
1:A:115:MET:SD	1:A:123:LYS:HB3	2.50	0.52
1:B:34:GLN:HE21	1:B:46:GLN:HG2	1.75	0.52
1:B:113:GLU:HB2	1:B:166:ARG:HB2	1.91	0.51
1:B:167:ILE:O	1:B:167:ILE:HG22	2.09	0.51
1:A:181:PRO:HA	1:A:184:VAL:CG2	2.40	0.51
1:B:114:HIS:HB2	1:B:165:LEU:CD1	2.40	0.51
1:A:108:ILE:HD12	1:A:132:ALA:HA	1.92	0.50
1:B:126:TRP:CZ3	1:B:147:LYS:HB2	2.46	0.50
1:B:150:VAL:HG12	1:B:151:LEU:N	2.27	0.49
1:A:182:GLY:O	1:A:183:GLU:CB	2.61	0.49
1:A:115:MET:SD	1:A:123:LYS:CB	3.00	0.49
1:B:104:ALA:N	1:B:106:THR:HG23	2.27	0.49
1:A:183:GLU:HG2	1:B:224:ASP:O	2.12	0.49
1:A:94:ALA:HB2	3:A:403:HOH:O	2.13	0.49
1:B:65:LYS:HB2	1:B:153:MET:CE	2.39	0.48
1:B:150:VAL:CG1	1:B:151:LEU:N	2.76	0.48
1:B:89:LEU:CB	1:B:90:PRO:CD	2.92	0.47
1:B:114:HIS:HB2	1:B:165:LEU:HD12	1.94	0.47
1:A:166:ARG:HB3	1:A:168:MET:CE	2.44	0.47
1:B:110:LYS:HD3	1:B:167:ILE:HG22	1.96	0.47
1:B:135:PRO:O	1:B:137:MET:N	2.48	0.47
1:B:113:GLU:HB3	1:B:125:GLU:OE2	2.15	0.47
1:B:112:VAL:HG12	1:B:113:GLU:N	2.30	0.47
1:A:197:LYS:CG	1:A:203:ARG:HG3	2.45	0.46
1:B:127:ARG:O	1:B:146:GLU:HG2	2.15	0.46
1:B:88:VAL:HG12	1:B:89:LEU:N	2.31	0.46
1:B:164:ASP:O	1:B:165:LEU:HD13	2.16	0.46
1:A:185:VAL:HG13	1:B:73:TYR:OH	2.16	0.45
1:B:28:ILE:O	1:B:29:VAL:C	2.55	0.45
1:B:143:ILE:HD12	1:B:156:LEU:HD11	1.98	0.45
1:B:166:ARG:O	1:B:168:MET:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:GLN:HA	1:A:47:TRP:O	2.17	0.45
1:A:197:LYS:HD2	1:A:203:ARG:CG	2.47	0.45
1:B:156:LEU:O	1:B:165:LEU:HD21	2.17	0.45
1:A:85:ALA:O	1:A:87:GLU:HG3	2.17	0.45
1:B:114:HIS:HB3	1:B:165:LEU:HD12	1.99	0.44
1:A:197:LYS:HG3	1:A:203:ARG:HG3	1.99	0.44
1:A:27:ASN:ND2	1:A:29:VAL:H	2.16	0.44
1:B:32:ARG:HH22	1:B:67:ASP:CG	2.21	0.44
1:B:233:LEU:O	1:B:234:VAL:HG23	2.18	0.43
1:B:166:ARG:CB	1:B:168:MET:HE3	2.48	0.43
1:B:166:ARG:HB2	1:B:168:MET:HE3	2.00	0.43
1:A:181:PRO:HA	1:A:184:VAL:HG21	2.00	0.43
1:B:161:LYS:C	1:B:163:GLY:N	2.70	0.43
1:B:53:ASP:HB3	1:B:63:LEU:HB2	2.01	0.43
1:B:28:ILE:O	1:B:31:CYS:HB2	2.18	0.43
1:B:113:GLU:HB2	1:B:168:MET:HE3	2.01	0.43
1:A:114:HIS:HE1	1:A:159:ASP:OD2	2.02	0.42
1:A:193:VAL:CG1	1:A:207:VAL:HG23	2.50	0.42
1:A:194:GLU:HA	1:A:203:ARG:O	2.20	0.42
1:B:167:ILE:O	1:B:168:MET:C	2.57	0.41
1:A:115:MET:SD	1:A:123:LYS:HB2	2.60	0.41
1:B:143:ILE:HG23	1:B:152:TYR:HB2	2.03	0.41
1:B:127:ARG:NE	1:B:146:GLU:HG3	2.35	0.41
1:B:112:VAL:CG1	1:B:113:GLU:N	2.84	0.41
1:B:28:ILE:HG22	1:B:28:ILE:O	2.20	0.41
1:A:131:LEU:O	1:A:132:ALA:HB2	2.20	0.41
1:B:146:GLU:H	1:B:146:GLU:HG2	1.61	0.41
1:A:119:GLU:H	1:A:119:GLU:HG2	1.66	0.41
1:B:54:GLN:O	1:B:56:PRO:HD3	2.22	0.40

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	197/242 (81%)	186 (94%)	9 (5%)	2 (1%)	19	16
1	B	164/242 (68%)	149 (91%)	13 (8%)	2 (1%)	16	12
All	All	361/484 (75%)	335 (93%)	22 (6%)	4 (1%)	17	14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	GLU
1	B	108	ILE
1	A	183	GLU
1	B	136	VAL

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	177/214 (83%)	164 (93%)	13 (7%)	17	18
1	B	152/214 (71%)	134 (88%)	18 (12%)	6	5
All	All	329/428 (77%)	298 (91%)	31 (9%)	11	10

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	26	ARG
1	A	58	ASN
1	A	97	ARG
1	A	114	HIS
1	A	118	THR
1	A	123	LYS
1	A	138	ASN
1	A	170	ASP
1	A	181	PRO
1	A	195	TYR
1	A	214	LYS

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Mol	Chain	Res	Type
1	A	226	PHE
1	B	32	ARG
1	B	41	ASN
1	B	50	THR
1	B	52	LEU
1	B	58	ASN
1	B	84	SER
1	B	89	LEU
1	B	106	THR
1	B	107	MET
1	B	114	HIS
1	B	125	GLU
1	B	161	LYS
1	B	165	LEU
1	B	168	MET
1	B	184	VAL
1	B	199	ASP
1	B	204	THR
1	B	226	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	41	ASN
1	A	58	ASN
1	A	114	HIS
1	A	138	ASN
1	A	210	GLN
1	B	34	GLN
1	B	46	GLN
1	B	58	ASN
1	B	155	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 ligands 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	PO4	A	301	-	4,4,4	1.42	0	6,6,6	0.27	0
2	PO4	A	302	-	4,4,4	1.20	0	6,6,6	0.27	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	PO4	A	301	-	-	0/0/0/0	0/0/0/0
2	PO4	A	302	-	-	0/0/0/0	0/0/0/0

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

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	201/242 (83%)	0.36	13 (6%) 22 22	21, 31, 68, 87	0
1	B	172/242 (71%)	1.30	52 (30%) 1 0	36, 60, 95, 106	0
All	All	373/484 (77%)	0.79	65 (17%) 2 2	21, 45, 90, 106	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	TYR	6.8
1	B	138	ASN	6.7
1	B	166	ARG	6.6
1	A	121	GLY	5.8
1	A	120	ASP	5.7
1	B	200	GLY	5.6
1	B	89	LEU	5.5
1	B	41	ASN	5.5
1	B	106	THR	5.4
1	B	165	LEU	5.2
1	B	167	ILE	4.9
1	B	163	GLY	4.9
1	B	156	LEU	4.8
1	B	157	LEU	4.8
1	B	196	ALA	4.7
1	B	115	MET	4.7
1	B	168	MET	4.7
1	B	111	ALA	4.6
1	B	133	ARG	4.6
1	B	141	PHE	4.5
1	B	161	LYS	4.3
1	B	146	GLU	4.0
1	A	122	SER	4.0
1	B	199	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	107	MET	3.9
1	A	183	GLU	3.8
1	B	136	VAL	3.7
1	B	113	GLU	3.7
1	B	112	VAL	3.7
1	A	182	GLY	3.6
1	B	164	ASP	3.5
1	A	123	LYS	3.5
1	B	126	TRP	3.4
1	B	127	ARG	3.4
1	B	197	LYS	3.4
1	B	110	LYS	3.2
1	B	162	GLU	3.2
1	A	235	LYS	3.2
1	B	114	HIS	3.2
1	B	147	LYS	3.1
1	B	139	THR	2.9
1	B	109	GLY	2.9
1	B	137	MET	2.8
1	B	90	PRO	2.8
1	A	41	ASN	2.8
1	A	170	ASP	2.7
1	B	201	SER	2.7
1	B	28	ILE	2.7
1	A	184	VAL	2.6
1	B	42	GLY	2.6
1	B	59	PRO	2.6
1	A	124	ASP	2.6
1	B	31	CYS	2.6
1	B	202	LYS	2.4
1	B	105	ASP	2.3
1	B	104	ALA	2.3
1	B	27	ASN	2.2
1	A	118	THR	2.2
1	B	140	TRP	2.2
1	A	119	GLU	2.1
1	B	158	ASP	2.1
1	B	184	VAL	2.1
1	B	134	ALA	2.1
1	B	154	TYR	2.0
1	B	40	GLY	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(\AA^2)	Q<0.9
2	PO4	A	301	5/5	0.99	0.18	2.40	30,32,34,36	0
2	PO4	A	302	5/5	0.97	0.14	-0.84	37,39,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.