



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 16, 2023 – 10:57 AM EST

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

<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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

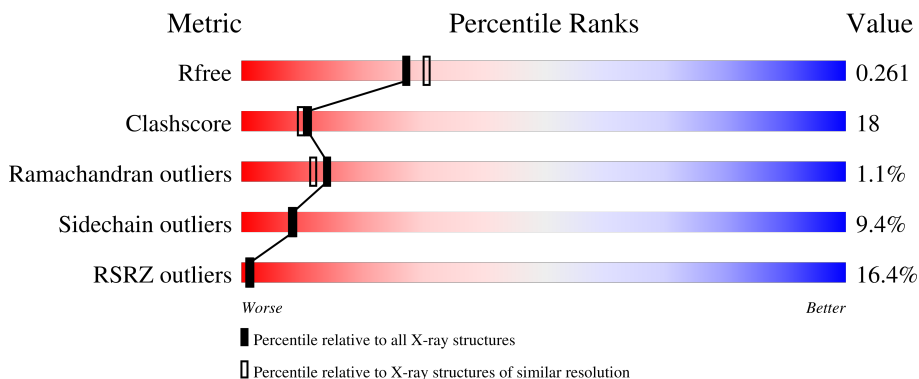
1 Overall quality at a glance

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}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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 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	242	<p>5% (poor fit), 63% (0 outliers), 17% (1 outlier), 17% (2+ outliers)</p>
1	B	242	<p>20% (poor fit), 42% (0 outliers), 25% (1 outlier), 29% (2+ outliers)</p>

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
			Total	C	N	O	S			
1	A	201	1626	1039	270	308	9	0	0	0
1	B	172	1397	903	225	260	9	0	0	0

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 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0

- Molecule 3 is water.

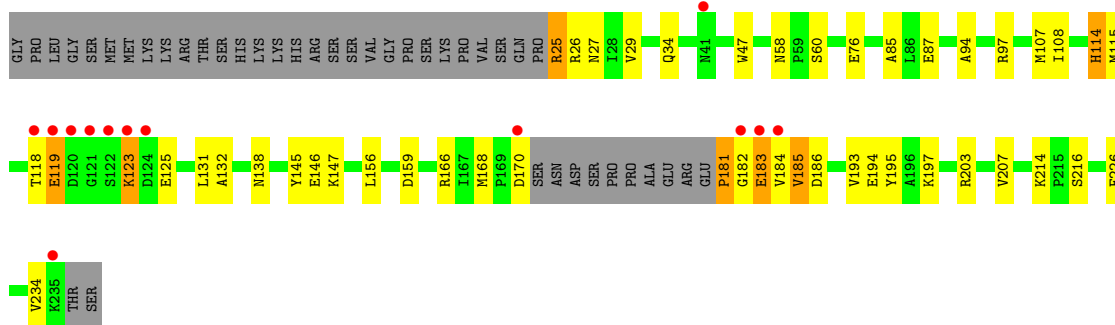
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	170	Total O 170 170	0	0
3	B	71	Total O 71 71	0	0

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.

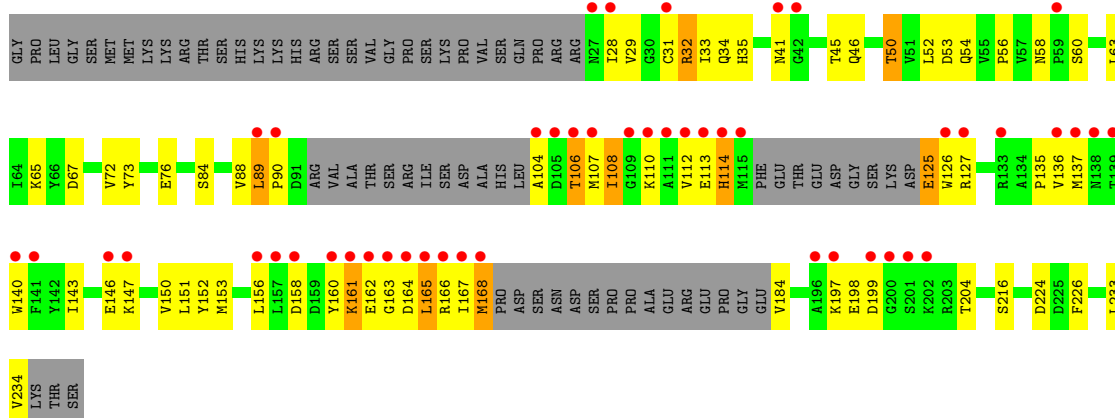
- Molecule 1: Spindlin-1

Chain A: 



- Molecule 1: Spindlin-1

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	40.79Å 84.88Å 136.57Å 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 2.12Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.225 , 0.260 0.226 , 0.261	Depositor DCC
R_{free} test set	2453 reflections (8.96%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtrriage
Anisotropy	0.079	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 69.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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, 0.333 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: 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:HG2	1:A:25:ARG:NH1	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:60:SER:O	1:B:76:GLU:HA	2.04	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: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:94:ALA:HB2	3:A:403:HOH:O	2.13	0.49
1:A:183:GLU:HG2	1:B:224:ASP:O	2.12	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:A:166:ARG:HB3	1:A:168:MET:CE	2.44	0.47
1:B:114:HIS:HB2	1:B:165:LEU:HD12	1.94	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:27:ASN:ND2	1:A:29:VAL:H	2.16	0.44
1:A:197:LYS:HG3	1:A:203:ARG:HG3	1.99	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:A:181:PRO:HA	1:A:184:VAL:HG21	2.00	0.43
1:B:166:ARG:HB2	1:B:168:MET:HE3	2.00	0.43
1:B:53:ASP:HB3	1:B:63:LEU:HB2	2.01	0.43
1:B:161:LYS:C	1:B:163:GLY:N	2.70	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:112:VAL:CG1	1:B:113:GLU:N	2.84	0.41
1:B:127:ARG:NE	1:B:146:GLU:HG3	2.35	0.41
1:B:143:ILE:HG23	1:B:152:TYR:HB2	2.03	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:A:119:GLU:H	1:A:119:GLU:HG2	1.66	0.41
1:B:146:GLU:HG2	1:B:146:GLU:H	1.61	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%)	15	14
1	B	164/242 (68%)	149 (91%)	13 (8%)	2 (1%)	13	10
All	All	361/484 (75%)	335 (93%)	22 (6%)	4 (1%)	14	12

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 [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	177/214 (83%)	164 (93%)	13 (7%)	14	15
1	B	152/214 (71%)	134 (88%)	18 (12%)	5	4
All	All	329/428 (77%)	298 (91%)	31 (9%)	8	8

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

Sometimes 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 [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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	302	-	4,4,4	1.74	1 (25%)	6,6,6	0.43	0
2	PO4	A	301	-	4,4,4	2.02	2 (50%)	6,6,6	0.45	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	PO4	P-O2	-2.62	1.46	1.54
2	A	301	PO4	P-O4	-2.22	1.47	1.54
2	A	302	PO4	P-O4	-2.06	1.48	1.54

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

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%)	0.29	13 (6%) 18 17	21, 31, 68, 87	0
1	B	172/242 (71%)	1.22	48 (27%) 0 0	36, 60, 95, 106	0
All	All	373/484 (77%)	0.72	61 (16%) 1 1	21, 45, 90, 106	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	TYR	6.7
1	B	138	ASN	6.5
1	B	166	ARG	6.4
1	A	121	GLY	5.6
1	A	120	ASP	5.5
1	B	200	GLY	5.4
1	B	41	ASN	5.3
1	B	89	LEU	5.3
1	B	106	THR	5.2
1	B	165	LEU	5.0
1	B	167	ILE	4.8
1	B	163	GLY	4.7
1	B	156	LEU	4.7
1	B	157	LEU	4.6
1	B	196	ALA	4.5
1	B	115	MET	4.5
1	B	168	MET	4.5
1	B	133	ARG	4.5
1	B	111	ALA	4.4
1	B	141	PHE	4.4
1	B	161	LYS	4.2
1	B	146	GLU	3.9
1	A	122	SER	3.8
1	B	199	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	107	MET	3.7
1	A	183	GLU	3.7
1	B	113	GLU	3.6
1	B	136	VAL	3.6
1	B	112	VAL	3.5
1	A	182	GLY	3.4
1	A	123	LYS	3.4
1	B	164	ASP	3.4
1	B	197	LYS	3.3
1	B	127	ARG	3.3
1	B	126	TRP	3.2
1	B	110	LYS	3.1
1	A	235	LYS	3.1
1	B	162	GLU	3.1
1	B	114	HIS	3.0
1	B	147	LYS	3.0
1	B	139	THR	2.8
1	B	109	GLY	2.8
1	B	137	MET	2.7
1	B	90	PRO	2.7
1	A	41	ASN	2.7
1	A	170	ASP	2.6
1	B	201	SER	2.6
1	B	28	ILE	2.6
1	B	42	GLY	2.5
1	A	184	VAL	2.5
1	A	124	ASP	2.5
1	B	59	PRO	2.5
1	B	31	CYS	2.4
1	B	202	LYS	2.4
1	B	105	ASP	2.2
1	B	104	ALA	2.1
1	B	27	ASN	2.1
1	A	118	THR	2.1
1	A	119	GLU	2.1
1	B	140	TRP	2.1
1	B	158	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

There are no monosaccharides 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
2	PO4	A	302	5/5	0.97	0.14	37,39,41,41	0
2	PO4	A	301	5/5	0.99	0.18	30,32,34,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.