



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:54 AM UTC

PDB ID : 1SKN / pdb_00001skn
Title : THE BINDING DOMAIN OF SKN-1 IN COMPLEX WITH DNA: A NEW
DNA-BINDING MOTIF
Authors : Rupert, P.B.; Daughdrill, G.W.; Bowerman, B.; Matthews, B.W.
Deposited on : 1998-03-30
Resolution : 2.50 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

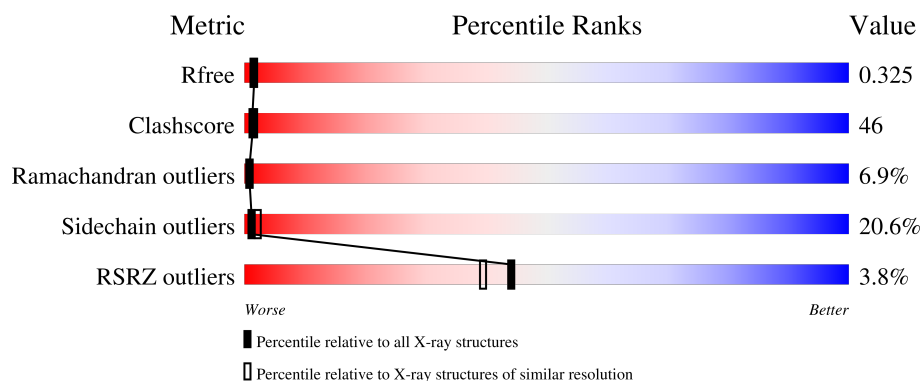
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

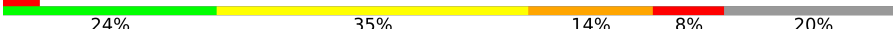
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	15	
2	B	15	
3	P	92	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 1263 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 DNA chain called DNA (5'-D(*TP*GP*AP*CP*AP*AP*TP*GP*TP*CP*AP*TP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	15	Total	C	N	O	P	0	0	0
			300	145	53	88	14			

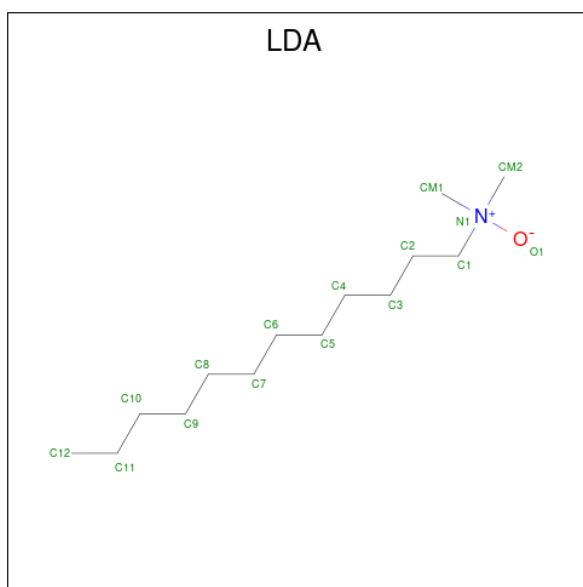
- Molecule 2 is a DNA chain called DNA (5'-D(*CP*AP*GP*GP*GP*AP*TP*GP*AP*CP*AP*TP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	15	Total	C	N	O	P	0	0	0
			309	148	59	88	14			

- Molecule 3 is a protein called DNA-BINDING DOMAIN OF SKN-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	74	Total	C	N	O	S	0	0	0
			610	365	126	116	3			

- Molecule 4 is LAURYL DIMETHYLAMINE-N-OXIDE (CCD ID: LDA) (formula: C₁₄H₃₁NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	P	1	Total	C	N	O	0	0
			16	14	1	1		

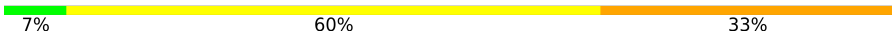
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	7	Total	O	0	0
			7	7		
5	B	4	Total	O	0	0
			4	4		
5	P	17	Total	O	0	0
			17	17		

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.

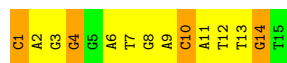
- Molecule 1: DNA (5'-D(*TP*GP*AP*CP*AP*AP*TP*GP*TP*CP*AP*TP*CP*CP*C)-3')

Chain A: 



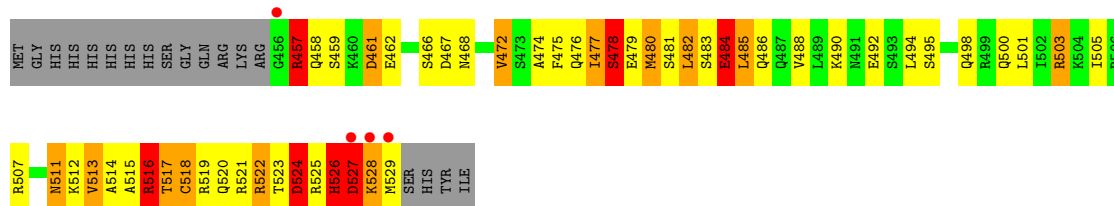
- Molecule 2: DNA (5'-D(*CP*AP*GP*GP*GP*AP*TP*GP*AP*CP*AP*TP*TP*GP*T)-3')

Chain B: 



- Molecule 3: DNA-BINDING DOMAIN OF SKN-1

Chain P: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	82.13Å 93.39Å 53.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.5 (20.00-2.50) 97.7 (20.00-2.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.19Å)	Xtriage
Refinement program	TNT, XTALVIEW	Depositor
R, R_{free}	0.220 , 0.337 0.227 , 0.325	Depositor DCC
R_{free} test set	1104 reflections (10.50%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 106.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	1263	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.15% 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: LDA

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	0.77	0/335	1.55	8/514 (1.6%)
2	B	0.80	0/347	1.54	4/535 (0.7%)
3	P	1.61	5/613 (0.8%)	2.19	23/813 (2.8%)
All	All	1.24	5/1295 (0.4%)	1.85	35/1862 (1.9%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	472	VAL	CA-C	-6.35	1.46	1.52
3	P	513	VAL	CA-CB	-5.80	1.47	1.54
3	P	484	GLU	N-CA	-5.28	1.39	1.46
3	P	483	SER	C-O	-5.15	1.17	1.24
3	P	461	ASP	CG-OD2	5.01	1.34	1.25

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	DC	P-O3'-C3'	-9.05	106.63	120.20
3	P	511	ASN	CA-C-N	-8.86	105.62	120.68
3	P	511	ASN	C-N-CA	-8.86	105.62	120.68
3	P	472	VAL	CA-C-O	-8.53	113.28	121.64
3	P	467	ASP	CA-CB-CG	-8.41	104.19	112.60
1	A	3	DA	P-O3'-C3'	8.37	132.75	120.20
3	P	524	ASP	CA-CB-CG	-7.62	104.98	112.60
3	P	475	PHE	CA-CB-CG	-7.06	106.74	113.80
1	A	11	DA	P-O5'-C5'	-6.93	109.60	120.00
3	P	468	ASN	CA-CB-CG	-6.90	105.70	112.60
1	A	5	DA	P-O5'-C5'	-6.63	110.06	120.00
3	P	527	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	4	DC	O3'-P-O5'	-6.46	94.30	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	14	DG	P-O5'-C5'	-6.41	110.38	120.00
3	P	485	LEU	CA-C-O	6.10	127.22	120.63
3	P	526	HIS	N-CA-CB	6.01	120.65	110.49
1	A	8	DG	O3'-P-O5'	-5.72	95.41	104.00
1	A	10	DC	P-O3'-C3'	5.72	128.78	120.20
1	A	8	DG	P-O5'-C5'	-5.63	111.56	120.00
3	P	477	ILE	N-CA-CB	-5.59	104.75	110.62
3	P	478	SER	N-CA-C	5.57	118.28	111.82
1	A	6	DA	O3'-P-O5'	-5.51	95.74	104.00
3	P	488	VAL	CA-C-O	5.48	126.75	121.05
3	P	472	VAL	O-C-N	5.46	129.00	122.95
3	P	500	GLN	N-CA-CB	-5.42	102.16	110.01
3	P	522	ARG	O-C-N	5.35	127.80	122.08
2	B	10	DC	C3'-C2'-C1'	-5.25	93.72	101.60
3	P	513	VAL	CA-C-N	-5.20	113.37	120.44
3	P	513	VAL	C-N-CA	-5.20	113.37	120.44
2	B	4	DG	O3'-P-O5'	-5.15	96.28	104.00
3	P	513	VAL	CA-C-O	5.12	126.28	120.95
3	P	500	GLN	CB-CA-C	5.07	118.84	110.88
3	P	482	LEU	CA-C-N	-5.06	112.27	120.72
3	P	482	LEU	C-N-CA	-5.06	112.27	120.72
3	P	461	ASP	N-CA-C	-5.01	105.95	111.71

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	300	0	171	18	0
2	B	309	0	171	28	0
3	P	610	0	626	59	0
4	P	16	0	31	3	0
5	A	7	0	0	2	0
5	B	4	0	0	0	0
5	P	17	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	1263	0	999	98	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 46.

All (98) 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:12:DT:H2''	1:A:13:DC:H5'	1.31	1.09
2:B:9:DA:C8	2:B:9:DA:H5'	2.06	0.89
2:B:1:DC:H2''	2:B:2:DA:H5'	1.59	0.81
1:A:12:DT:C2'	1:A:13:DC:H5'	2.11	0.79
2:B:9:DA:H5'	2:B:9:DA:H8	1.47	0.75
1:A:4:DC:H1'	1:A:5:DA:H5'	1.70	0.74
1:A:12:DT:H2''	1:A:13:DC:C5'	2.17	0.73
3:P:526:HIS:O	3:P:527:ASP:C	2.31	0.71
2:B:8:DG:H2''	2:B:9:DA:C5'	2.22	0.70
2:B:6:DA:H2''	2:B:7:DT:C5'	2.22	0.70
3:P:476:GLN:O	3:P:480:MET:HB3	1.92	0.69
1:A:8:DG:O6	3:P:519:ARG:NH2	2.26	0.69
3:P:492:GLU:OE2	4:P:16:LDA:HM21	1.93	0.69
3:P:478:SER:HB2	3:P:479:GLU:OE2	1.92	0.68
3:P:516:ARG:O	3:P:517:THR:C	2.35	0.68
3:P:517:THR:O	3:P:518:CYS:C	2.36	0.68
3:P:526:HIS:O	3:P:528:LYS:N	2.26	0.67
3:P:520:GLN:O	3:P:524:ASP:N	2.29	0.65
3:P:512:LYS:HG3	3:P:516:ARG:NH1	2.12	0.65
3:P:517:THR:O	3:P:520:GLN:N	2.29	0.65
2:B:12:DT:H1'	2:B:13:DT:H5''	1.77	0.64
2:B:6:DA:H2''	2:B:7:DT:H5'	1.79	0.64
2:B:2:DA:H2'	2:B:3:DG:C8	2.33	0.64
2:B:7:DT:OP2	3:P:521:ARG:NH2	2.29	0.63
2:B:12:DT:H1'	2:B:13:DT:C5'	2.29	0.63
3:P:518:CYS:N	3:P:521:ARG:NH1	2.46	0.62
2:B:6:DA:H2''	2:B:7:DT:O5'	1.98	0.62
3:P:482:LEU:O	3:P:486:GLN:HG3	1.99	0.62
3:P:518:CYS:HA	3:P:521:ARG:CZ	2.29	0.62
3:P:462:GLU:CD	3:P:474:ALA:HB3	2.25	0.61
3:P:518:CYS:N	3:P:521:ARG:HH12	1.98	0.61
3:P:457:ARG:HD3	3:P:458:GLN:O	2.00	0.61
3:P:520:GLN:O	3:P:524:ASP:HB2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:P:494:LEU:O	3:P:495:SER:C	2.41	0.60
1:A:8:DG:OP2	3:P:516:ARG:NH1	2.34	0.60
2:B:8:DG:H2''	2:B:9:DA:H5'	1.84	0.59
2:B:1:DC:C6	2:B:1:DC:H5'	2.37	0.59
3:P:517:THR:OG1	3:P:521:ARG:NH1	2.29	0.59
2:B:2:DA:H2''	2:B:3:DG:O5'	2.01	0.59
3:P:492:GLU:CD	4:P:16:LDA:HM21	2.27	0.58
1:A:13:DC:H1'	1:A:14:DC:H5'	1.86	0.58
2:B:4:DG:OP2	3:P:503:ARG:NH2	2.36	0.57
3:P:518:CYS:HA	3:P:521:ARG:NH2	2.18	0.57
3:P:501:LEU:O	3:P:505:ILE:HG13	2.04	0.57
1:A:7:DT:H3'	5:A:117:HOH:O	2.04	0.56
2:B:1:DC:C2'	2:B:2:DA:H5'	2.33	0.54
1:A:12:DT:C4	1:A:13:DC:N4	2.76	0.54
3:P:518:CYS:O	3:P:522:ARG:N	2.35	0.54
1:A:6:DA:H2''	1:A:7:DT:H5'	1.90	0.54
3:P:514:ALA:O	3:P:515:ALA:C	2.52	0.52
3:P:457:ARG:HG2	3:P:458:GLN:H	1.74	0.52
3:P:492:GLU:OE1	4:P:16:LDA:HM21	2.09	0.52
3:P:503:ARG:O	3:P:503:ARG:NH1	2.44	0.51
3:P:515:ALA:O	3:P:516:ARG:O	2.29	0.51
3:P:481:SER:O	3:P:484:GLU:N	2.45	0.50
2:B:6:DA:C2'	2:B:7:DT:H5'	2.40	0.49
2:B:7:DT:H2'	3:P:518:CYS:SG	2.51	0.49
3:P:523:THR:O	3:P:527:ASP:HB3	2.12	0.49
2:B:1:DC:H6	2:B:1:DC:C5'	2.25	0.49
1:A:14:DC:H2''	1:A:15:DC:OP2	2.13	0.49
3:P:521:ARG:O	3:P:525:ARG:HG3	2.12	0.48
3:P:472:VAL:HG23	3:P:477:ILE:HG13	1.95	0.48
2:B:10:DC:H2''	2:B:11:DA:OP2	2.13	0.48
3:P:517:THR:CB	3:P:521:ARG:HH12	2.25	0.48
1:A:5:DA:C2	2:B:14:DG:C2	3.02	0.48
3:P:521:ARG:NH1	3:P:521:ARG:HG3	2.29	0.47
2:B:1:DC:C6	2:B:1:DC:C5'	2.98	0.47
3:P:459:SER:OG	3:P:462:GLU:HG2	2.15	0.46
3:P:501:LEU:HD12	3:P:501:LEU:HA	1.67	0.46
3:P:511:ASN:O	3:P:512:LYS:C	2.52	0.46
3:P:512:LYS:HG3	3:P:516:ARG:HH12	1.78	0.46
3:P:521:ARG:HG3	3:P:521:ARG:HH11	1.81	0.45
3:P:515:ALA:O	3:P:516:ARG:C	2.59	0.45
2:B:8:DG:C6	2:B:9:DA:C6	3.05	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2:DA:H2''	2:B:3:DG:C5'	2.46	0.45
3:P:459:SER:OG	3:P:462:GLU:N	2.38	0.44
1:A:1:DT:H2''	1:A:2:DG:C8	2.53	0.43
3:P:519:ARG:NH1	5:P:107:HOH:O	2.31	0.43
1:A:12:DT:N3	1:A:13:DC:C4	2.86	0.43
3:P:472:VAL:HG23	3:P:477:ILE:CG1	2.48	0.43
1:A:5:DA:H2''	1:A:6:DA:OP2	2.18	0.43
3:P:495:SER:N	3:P:498:GLN:OE1	2.52	0.43
3:P:481:SER:O	3:P:484:GLU:HB2	2.19	0.43
3:P:517:THR:CB	3:P:521:ARG:NH1	2.81	0.43
2:B:11:DA:H1'	2:B:12:DT:H5'	2.01	0.42
2:B:12:DT:H1'	2:B:13:DT:H5'	1.99	0.42
3:P:507:ARG:NH2	5:P:118:HOH:O	2.53	0.42
3:P:523:THR:O	3:P:523:THR:HG22	2.19	0.42
5:A:108:HOH:O	3:P:461:ASP:HB2	2.20	0.41
1:A:5:DA:O5'	1:A:5:DA:H2'	2.20	0.41
3:P:477:ILE:HG21	3:P:477:ILE:HD13	1.79	0.41
3:P:494:LEU:HA	3:P:498:GLN:OE1	2.20	0.41
1:A:1:DT:H4'	1:A:2:DG:OP1	2.20	0.41
1:A:2:DG:C2	1:A:3:DA:C6	3.09	0.41
2:B:12:DT:H2''	2:B:13:DT:OP2	2.21	0.41
2:B:7:DT:C2'	3:P:518:CYS:SG	3.10	0.40
3:P:480:MET:SD	3:P:485:LEU:HA	2.62	0.40
3:P:517:THR:C	3:P:521:ARG:NH1	2.79	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
3	P	72/92 (78%)	60 (83%)	7 (10%)	5 (7%)	1 1

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	P	516	ARG
3	P	527	ASP
3	P	517	THR
3	P	518	CYS
3	P	457	ARG

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
3	P	68/84 (81%)	54 (79%)	14 (21%)	1 2

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	P	457	ARG
3	P	466	SER
3	P	478	SER
3	P	480	MET
3	P	484	GLU
3	P	490	LYS
3	P	503	ARG
3	P	513	VAL
3	P	516	ARG
3	P	524	ASP
3	P	526	HIS
3	P	527	ASP
3	P	528	LYS
3	P	529	MET

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

Mol	Chain	Res	Type
3	P	463	GLN

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Mol	Chain	Res	Type
3	P	468	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
4	LDA	P	16	-	13,15,15	0.98	1 (7%)	14,17,17	0.69	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
4	LDA	P	16	-	-	8/13/13/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	16	LDA	C1-N1	-3.41	1.48	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	P	16	LDA	C2-C1-N1-O1
4	P	16	LDA	C2-C1-N1-CM1
4	P	16	LDA	C2-C1-N1-CM2
4	P	16	LDA	C11-C10-C9-C8
4	P	16	LDA	C5-C6-C7-C8
4	P	16	LDA	C4-C5-C6-C7
4	P	16	LDA	C9-C10-C11-C12
4	P	16	LDA	C3-C4-C5-C6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	16	LDA	3	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	15/15 (100%)	0.35	0	100 100	44, 58, 86, 91	0
2	B	15/15 (100%)	0.18	0	100 100	42, 60, 72, 73	0
3	P	74/92 (80%)	0.10	4 (5%)	31 27	24, 40, 98, 100	0
All	All	104/122 (85%)	0.15	4 (3%)	44 39	24, 47, 96, 100	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	P	527	ASP	2.9
3	P	528	LYS	2.8
3	P	456	GLY	2.5
3	P	529	MET	2.1

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
4	LDA	P	16	16/16	0.92	0.12	13,47,51,52	16

6.5 Other polymers [i](#)

There are no such residues in this entry.