



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 07:04 PM GMT

PDB ID : 3D3F
Title : Crystal Structure of Yvgn and cofactor NADPH from Bacillus subtilis
Authors : Zhou, Y.F.; Lei, J.; Su, X.D.
Deposited on : 2008-05-10
Resolution : 2.40 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

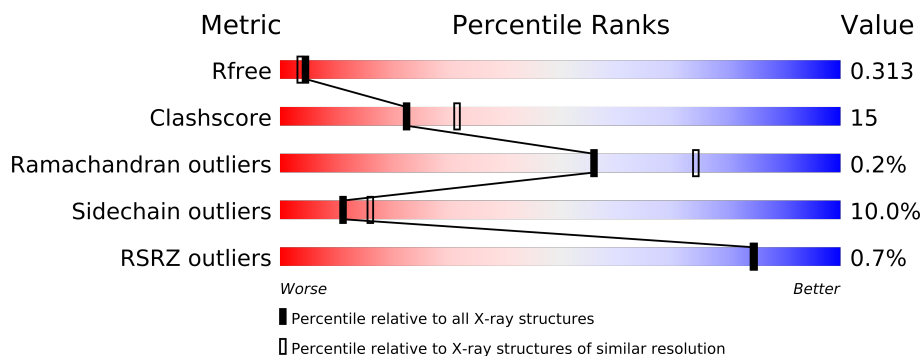
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.40 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}	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	275	
1	B	275	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 4629 atoms, of which 0 are hydrogen and 0 are deuterium.

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 YvgN protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	1	0
			2226	1417	378	426	5			
1	B	275	Total	C	N	O	S	0	1	0
			2226	1417	378	426	5			

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is water.

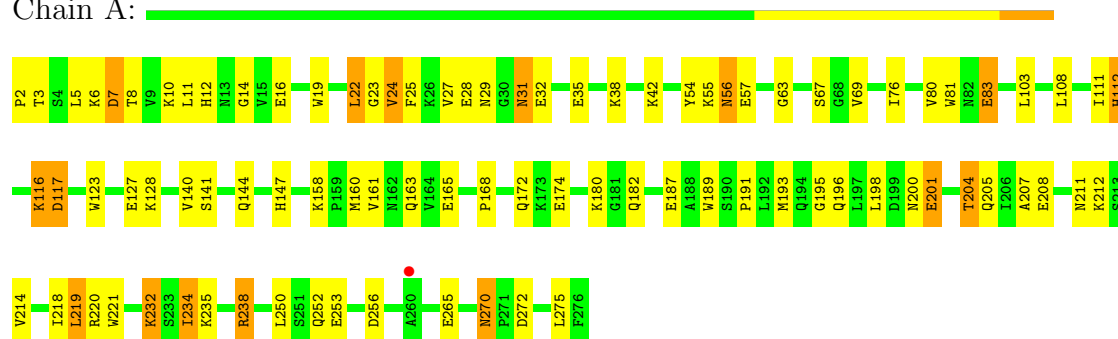
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total 41	O 41	0	0
3	B	40	Total 40	O 40	0	0

3 Residue-property plots

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 ($RSRZ > 2$). 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.

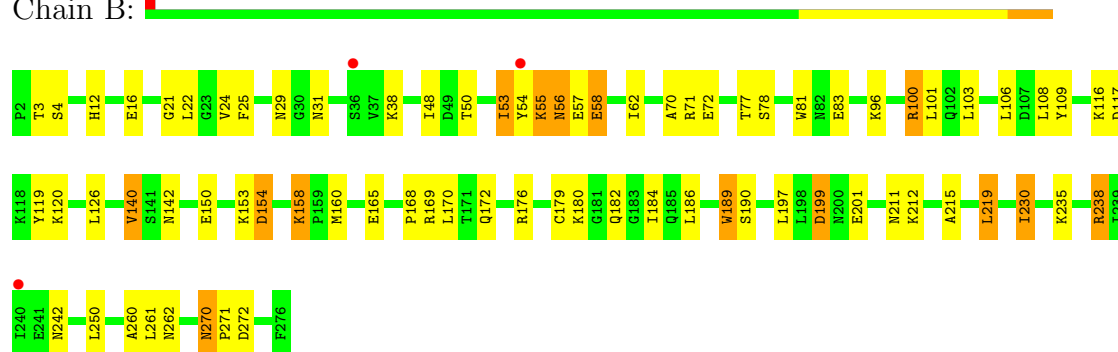
• Molecule 1: YvgN protein

Chain A:



• Molecule 1: YvgN protein

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.52Å 123.28Å 57.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 67.41 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.8 (20.00-2.40) 89.6 (67.41-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.236 , 0.306 0.250 , 0.313	Depositor DCC
R_{free} test set	1156 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.695	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 12.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Outliers	3 of 23303 reflections (0.013%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4629	wwPDB-VP
Average B, all atoms (Å ²)	9.0	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 37.44 % 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 4.2898e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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: NDP

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	1.08	5/2271 (0.2%)	0.98	6/3065 (0.2%)
1	B	1.00	2/2271 (0.1%)	0.95	5/3065 (0.2%)
All	All	1.04	7/4542 (0.2%)	0.97	11/6130 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	24	VAL	CB-CG1	-8.52	1.34	1.52
1	A	117	ASP	CB-CG	-8.06	1.34	1.51
1	A	57	GLU	CD-OE2	-6.96	1.18	1.25
1	A	57	GLU	CD-OE1	-6.44	1.18	1.25
1	A	24	VAL	CB-CG2	-5.49	1.41	1.52
1	B	189	TRP	CE3-CZ3	-5.14	1.29	1.38
1	B	16	GLU	CG-CD	-5.07	1.44	1.51

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	116	LYS	CB-CA-C	-7.24	95.92	110.40
1	A	116	LYS	N-CA-C	7.06	130.07	111.00
1	A	7	ASP	CB-CA-C	-5.85	98.69	110.40
1	B	260	ALA	CB-CA-C	5.55	118.43	110.10
1	A	29	ASN	CB-CA-C	5.24	120.88	110.40
1	B	219	LEU	CB-CG-CD1	5.16	119.77	111.00
1	A	22	LEU	CA-CB-CG	5.15	127.15	115.30
1	B	140	VAL	CG1-CB-CG2	5.13	119.12	110.90
1	A	220	ARG	NE-CZ-NH1	-5.07	117.76	120.30
1	A	22	LEU	CB-CG-CD2	5.05	119.59	111.00
1	B	116	LYS	C-N-CA	-5.05	109.08	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2226	0	2210	80	0
1	B	2226	0	2209	65	0
2	A	48	0	26	2	0
2	B	48	0	26	8	0
3	A	41	0	0	4	0
3	B	40	0	0	1	0
All	All	4629	0	4471	133	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (133) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:55:LYS:HG2	1:B:55:LYS:CG	1.76	1.15
1:A:8:THR:HG21	1:A:16:GLU:HB3	1.29	1.14
1:A:55:LYS:CG	1:B:55:LYS:HG2	1.81	1.09
1:A:8:THR:CG2	1:A:16:GLU:HB3	1.92	0.98
1:A:144:GLN:H	1:A:147:HIS:HD2	1.14	0.94
1:A:27:VAL:CG2	1:A:234:ILE:HD11	2.03	0.88
1:A:27:VAL:HG22	1:A:234:ILE:CD1	2.04	0.87
1:B:242:ASN:HD21	2:B:1:NDP:H2A	1.38	0.86
1:B:190:SER:HA	2:B:1:NDP:O1N	1.76	0.84
1:A:28:GLU:HG3	1:B:96:LYS:NZ	1.94	0.82
1:A:8:THR:HG21	1:A:16:GLU:CB	2.11	0.80
1:A:27:VAL:CG2	1:A:234:ILE:CD1	2.59	0.80
1:A:144:GLN:H	1:A:147:HIS:CD2	2.01	0.79
1:A:55:LYS:CE	1:B:55:LYS:HD3	2.12	0.78
1:A:200:ASN:O	1:A:204:THR:HG23	1.85	0.77
1:B:54:TYR:OH	2:B:1:NDP:H2D	1.85	0.77
1:B:158:LYS:HE3	1:B:182:GLN:O	1.86	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:195:GLY:HA3	3:A:296:HOH:O	1.88	0.72
1:B:270:ASN:ND2	1:B:272:ASP:H	1.88	0.71
1:A:55:LYS:HG2	1:B:55:LYS:HG2	0.84	0.71
1:A:28:GLU:HG3	1:B:96:LYS:HZ3	1.53	0.70
1:A:27:VAL:HG22	1:A:234:ILE:HD11	1.69	0.70
1:A:55:LYS:HE3	1:B:55:LYS:HD3	1.73	0.70
1:A:270:ASN:HD22	1:A:272:ASP:H	1.41	0.69
1:A:27:VAL:HG22	1:A:234:ILE:HD13	1.75	0.69
1:B:242:ASN:ND2	2:B:1:NDP:H2A	2.07	0.68
1:B:56:ASN:H	1:B:56:ASN:ND2	1.92	0.68
1:A:28:GLU:CG	1:B:96:LYS:NZ	2.58	0.67
1:B:53:ILE:HG12	1:B:81:TRP:CD2	2.33	0.64
1:A:27:VAL:HG21	1:A:234:ILE:HD11	1.78	0.64
1:B:158:LYS:CE	1:B:182:GLN:O	2.46	0.64
1:B:120:LYS:NZ	1:B:154:ASP:HB2	2.13	0.64
1:A:144:GLN:N	1:A:147:HIS:HD2	1.93	0.62
1:A:54:TYR:O	1:A:55:LYS:HB2	1.99	0.62
1:A:165:GLU:O	1:A:172:GLN:NE2	2.28	0.62
1:B:190:SER:CA	2:B:1:NDP:O1N	2.48	0.62
1:B:56:ASN:HD22	1:B:56:ASN:H	1.49	0.61
1:B:270:ASN:HD22	1:B:271:PRO:N	1.99	0.61
1:A:270:ASN:ND2	1:A:272:ASP:HB2	2.17	0.59
1:B:211:ASN:O	1:B:212:LYS:HD3	2.02	0.59
1:A:116:LYS:HD2	1:A:275:LEU:HB3	1.84	0.59
1:A:3:THR:N	1:A:7:ASP:OD2	2.36	0.59
1:A:8:THR:CG2	1:A:16:GLU:CB	2.75	0.58
1:A:27:VAL:HG21	1:A:234:ILE:CD1	2.34	0.57
1:A:12:HIS:CE1	1:A:160:MET:HG3	2.39	0.57
1:B:142:ASN:O	1:B:271:PRO:HB3	2.04	0.57
1:A:28:GLU:CG	1:B:96:LYS:HZ2	2.17	0.57
1:B:172:GLN:O	1:B:176:ARG:HG3	2.04	0.57
1:B:12:HIS:CE1	1:B:160:MET:HG2	2.40	0.57
1:B:168:PRO:O	1:B:262:ASN:HA	2.05	0.56
1:A:56:ASN:C	1:A:56:ASN:HD22	2.08	0.56
1:A:207:ALA:HB1	1:A:212:LYS:O	2.06	0.56
1:A:27:VAL:CG2	1:A:234:ILE:HD13	2.34	0.56
1:B:270:ASN:HD22	1:B:272:ASP:H	1.54	0.55
1:A:55:LYS:CD	1:B:55:LYS:HD3	2.36	0.55
1:A:198:LEU:HA	1:A:214:VAL:HG11	1.88	0.55
1:B:57:GLU:OE1	1:B:100:ARG:NH1	2.39	0.55
1:B:165:GLU:HB2	1:B:189:TRP:CZ2	2.42	0.55
1:A:270:ASN:HD21	1:A:272:ASP:HB2	1.72	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:53:ILE:HG12	1:B:81:TRP:CG	2.42	0.54
1:A:232:LYS:O	2:A:1:NDP:H8A	2.08	0.54
1:B:58:GLU:O	1:B:62:ILE:HG13	2.08	0.53
1:A:111:ILE:O	1:A:141:SER:HB3	2.09	0.53
1:B:165:GLU:HB2	1:B:189:TRP:CH2	2.44	0.53
1:B:50:THR:O	1:B:78:SER:HA	2.09	0.53
1:B:29:ASN:HD21	1:B:58:GLU:HG3	1.75	0.52
1:B:189:TRP:CD2	2:B:1:NDP:H41N	2.44	0.52
1:A:25:PHE:CZ	1:B:83:GLU:HG3	2.45	0.52
1:A:8:THR:HG23	1:A:16:GLU:HB3	1.86	0.52
1:B:165:GLU:O	1:B:172:GLN:NE2	2.38	0.52
1:A:158:LYS:HD3	1:A:182:GLN:O	2.10	0.51
1:A:117:ASP:O	3:A:277:HOH:O	2.19	0.51
2:B:1:NDP:H3D	2:B:1:NDP:H6N	1.93	0.51
1:A:235:LYS:HB2	1:A:238:ARG:HG3	1.94	0.50
1:B:96:LYS:O	1:B:100:ARG:HG2	2.11	0.50
1:A:76:ILE:HG21	1:A:103:LEU:HD11	1.94	0.50
1:A:11:LEU:HD13	1:A:160:MET:HB3	1.94	0.49
1:A:193:MET:CE	1:A:196:GLN:HB2	2.41	0.49
1:B:215:ALA:O	1:B:219:LEU:HG	2.13	0.49
1:A:32:GLU:HG2	1:A:234:ILE:HG13	1.94	0.48
1:B:21:GLY:HA3	1:B:230:ILE:HG22	1.96	0.48
1:A:81:TRP:CD1	1:A:83:GLU:HB3	2.49	0.48
1:B:270:ASN:HD22	1:B:271:PRO:CD	2.27	0.48
1:A:218:ILE:O	1:A:221:TRP:HB3	2.14	0.47
1:A:200:ASN:O	1:A:204:THR:CG2	2.61	0.47
1:B:270:ASN:HD22	1:B:270:ASN:C	2.18	0.47
1:A:211:ASN:O	1:A:212:LYS:HD3	2.15	0.46
1:A:270:ASN:ND2	1:A:272:ASP:H	2.12	0.46
1:A:201:GLU:H	1:A:201:GLU:HG2	1.60	0.46
1:B:25:PHE:HB2	2:B:1:NDP:H2D	1.98	0.46
1:B:109:TYR:HD2	1:B:126:LEU:HD22	1.81	0.46
1:A:27:VAL:HG23	1:A:232:LYS:HE2	1.98	0.45
1:B:120:LYS:HZ1	1:B:154:ASP:HB2	1.80	0.45
1:A:163:GLN:NE2	1:A:187:GLU:OE2	2.47	0.45
1:A:158:LYS:HB3	3:A:308:HOH:O	2.16	0.44
1:B:56:ASN:N	1:B:56:ASN:HD22	2.11	0.44
1:B:197:LEU:C	1:B:199:ASP:H	2.19	0.44
1:B:169:ARG:HG2	1:B:197:LEU:HD21	1.98	0.44
1:B:77:THR:HA	1:B:108:LEU:O	2.17	0.44
1:B:22:LEU:HB3	1:B:48:ILE:HD13	1.98	0.44
1:B:71:ARG:HD2	1:B:71:ARG:HA	1.85	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:2:PRO:HG3	1:A:19:TRP:CD1	2.54	0.43
1:B:4:SER:HA	3:B:315:HOH:O	2.17	0.43
1:A:55:LYS:HE3	1:B:55:LYS:CD	2.46	0.43
1:B:56:ASN:N	1:B:56:ASN:ND2	2.54	0.43
1:B:179:CYS:HB3	1:B:184:ILE:O	2.17	0.43
1:A:165:GLU:HB2	1:A:189:TRP:CH2	2.53	0.43
1:B:101:LEU:HB3	1:B:103:LEU:HG	1.99	0.43
1:A:35:GLU:OE2	1:A:38:LYS:HE2	2.17	0.43
1:B:70:ALA:O	1:B:71:ARG:C	2.58	0.43
1:A:63:GLY:O	1:A:67:SER:HB3	2.19	0.42
1:A:24:VAL:O	1:A:27:VAL:HB	2.19	0.42
1:A:198:LEU:HB3	3:A:302:HOH:O	2.19	0.42
1:B:81:TRP:HD1	1:B:83:GLU:HB3	1.85	0.42
1:A:193:MET:HE2	1:A:196:GLN:HB2	2.02	0.42
1:A:123:TRP:O	1:A:127:GLU:HG3	2.19	0.42
1:A:31:ASN:HA	1:A:31:ASN:HD22	1.57	0.42
1:B:165:GLU:O	1:B:170:LEU:HD23	2.20	0.42
1:A:28:GLU:HG2	1:B:96:LYS:HZ2	1.85	0.41
1:A:204:THR:O	1:A:208:GLU:HG3	2.20	0.41
1:B:106:LEU:HD12	1:B:109:TYR:CD1	2.55	0.41
1:A:108:LEU:HD21	1:A:161:VAL:HG21	2.02	0.41
1:A:219:LEU:HA	1:A:219:LEU:HD12	1.98	0.41
1:A:193:MET:HE3	1:A:196:GLN:HB2	2.02	0.41
1:A:252:GLN:NE2	1:A:256:ASP:OD1	2.53	0.41
1:A:10:LYS:HG2	1:A:14:GLY:HA2	2.02	0.41
1:A:189:TRP:CD2	2:A:1:NDP:H41N	2.56	0.41
1:B:119:TYR:OH	1:B:150:GLU:HG2	2.21	0.40
1:A:23:GLY:O	1:A:232:LYS:HB2	2.20	0.40
1:A:27:VAL:CG1	1:A:234:ILE:HD11	2.52	0.40
1:B:235:LYS:HD2	1:B:238:ARG:HE	1.86	0.40
1:A:80:VAL:O	1:A:112:HIS:HB2	2.21	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	273/275 (99%)	264 (97%)	9 (3%)	0	100	100
1	B	273/275 (99%)	248 (91%)	24 (9%)	1 (0%)	43	61
All	All	546/550 (99%)	512 (94%)	33 (6%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	261	LEU

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	241/241 (100%)	216 (90%)	25 (10%)	10	14
1	B	241/241 (100%)	218 (90%)	23 (10%)	12	18
All	All	482/482 (100%)	434 (90%)	48 (10%)	11	16

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	6	LYS
1	A	22	LEU
1	A	31	ASN
1	A	42	LYS
1	A	56	ASN
1	A	69	VAL
1	A	83	GLU
1	A	112	HIS
1	A	128	LYS
1	A	140	VAL
1	A	168	PRO
1	A	174	GLU
1	A	180	LYS
1	A	201	GLU
1	A	204	THR

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Mol	Chain	Res	Type
1	A	205	GLN
1	A	219	LEU
1	A	232	LYS
1	A	234	ILE
1	A	238	ARG
1	A	250	LEU
1	A	253	GLU
1	A	265	GLU
1	A	270	ASN
1	B	3	THR
1	B	24	VAL
1	B	31	ASN
1	B	38	LYS
1	B	53	ILE
1	B	55	LYS
1	B	56	ASN
1	B	58	GLU
1	B	72	GLU
1	B	100	ARG
1	B	117	ASP
1	B	140	VAL
1	B	153	LYS
1	B	154	ASP
1	B	158	LYS
1	B	180	LYS
1	B	186	LEU
1	B	199	ASP
1	B	201	GLU
1	B	230	ILE
1	B	238	ARG
1	B	250	LEU
1	B	270	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	31	ASN
1	A	56	ASN
1	A	147	HIS
1	A	162	ASN
1	A	252	GLN

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Mol	Chain	Res	Type
1	A	270	ASN
1	B	12	HIS
1	B	56	ASN
1	B	162	ASN
1	B	185	GLN
1	B	211	ASN
1	B	270	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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	NDP	A	1	-	52,52,52	1.44	7 (13%)	80,80,80	1.79	8 (10%)
2	NDP	B	1	-	52,52,52	1.46	5 (9%)	80,80,80	1.95	19 (23%)

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	NDP	A	1	-	-	0/35/77/77	0/3/5/5
2	NDP	B	1	-	-	0/35/77/77	0/3/5/5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	NDP	O7N-C7N	5.86	1.39	1.24
2	B	1	NDP	O7N-C7N	5.59	1.39	1.24
2	B	1	NDP	C2A-N3A	4.30	1.40	1.32
2	A	1	NDP	C2A-N3A	3.48	1.39	1.32
2	B	1	NDP	C2A-N1A	3.35	1.40	1.33
2	B	1	NDP	C8A-N9A	3.12	1.41	1.36
2	A	1	NDP	C2A-N1A	2.71	1.39	1.33
2	A	1	NDP	C2N-C3N	2.55	1.39	1.34
2	B	1	NDP	C2N-C3N	2.54	1.39	1.34
2	A	1	NDP	C6N-C5N	2.50	1.38	1.33
2	A	1	NDP	PN-O3	2.27	1.64	1.59
2	A	1	NDP	PA-O3	2.19	1.63	1.59

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NDP	N3A-C2A-N1A	-10.28	120.12	128.71
2	B	1	NDP	N3A-C2A-N1A	-7.07	122.80	128.71
2	B	1	NDP	O4B-C1B-N9A	6.27	114.27	108.44
2	A	1	NDP	O4B-C1B-N9A	5.68	113.73	108.44
2	B	1	NDP	N3A-C4A-N9A	5.34	135.07	125.43
2	B	1	NDP	C3B-C2B-C1B	-4.00	94.95	102.73
2	B	1	NDP	C8A-N9A-C1B	-3.84	118.81	126.38
2	B	1	NDP	C1B-N9A-C4A	3.83	133.26	126.64
2	A	1	NDP	C1D-N1N-C2N	-3.54	114.98	121.02
2	B	1	NDP	C4B-O4B-C1B	-3.31	106.16	109.75
2	B	1	NDP	O5D-PN-O1N	-3.13	97.11	109.37
2	A	1	NDP	N3A-C4A-N9A	3.13	131.08	125.43
2	B	1	NDP	O7N-C7N-C3N	-3.12	114.65	120.51
2	B	1	NDP	C5A-C4A-N3A	-2.82	119.56	125.70
2	B	1	NDP	C2B-C1B-N9A	2.79	124.30	113.74
2	B	1	NDP	C3N-C7N-N7N	2.68	122.63	117.48
2	B	1	NDP	N6A-C6A-N1A	2.63	124.53	119.36
2	B	1	NDP	C5A-C6A-N6A	-2.49	115.10	120.72
2	B	1	NDP	C6A-C5A-C4A	2.47	121.77	117.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	NDP	O2X-P2B-O2B	2.34	113.83	107.09
2	B	1	NDP	O3D-C3D-C2D	2.31	119.36	111.83
2	B	1	NDP	O2A-PA-O3	2.25	115.81	105.14
2	A	1	NDP	C5A-C4A-N3A	-2.23	120.84	125.70
2	B	1	NDP	C1D-N1N-C2N	-2.13	117.39	121.02
2	A	1	NDP	C4B-O4B-C1B	2.07	112.00	109.75
2	B	1	NDP	O3B-C3B-C4B	-2.05	105.05	111.08
2	A	1	NDP	C2D-C1D-N1N	-2.03	108.03	113.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	275/275 (100%)	-0.21	1 (0%) 90 90	4, 9, 16, 24	1 (0%)
1	B	275/275 (100%)	-0.12	3 (1%) 77 77	2, 9, 16, 22	1 (0%)
All	All	550/550 (100%)	-0.16	4 (0%) 84 84	2, 9, 16, 24	2 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	36[A]	SER	3.5
1	A	260	ALA	3.0
1	B	240	ILE	2.5
1	B	54	TYR	2.2

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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NDP	A	1	48/48	0.14	0.10	2,12,16,18	0
2	NDP	B	1	48/48	0.14	-0.01	5,11,16,19	0

6.5 Other polymers ⓘ

There are no such residues in this entry.