



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 05:58 AM GMT

PDB ID : 1NVT  
Title : Crystal structure of Shikimate Dehydrogenase (AROE or MJ1084) in complex with NADP+  
Authors : Padyana, A.K.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2003-02-04  
Resolution : 2.35 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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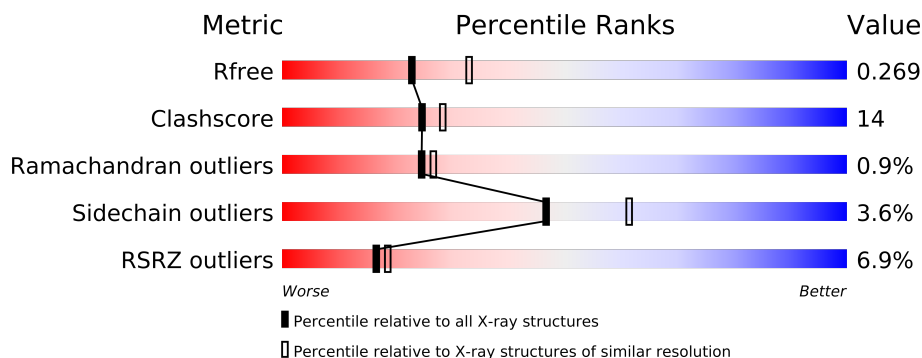
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.35 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	3327 (2.40-2.32)
Clashscore	79885	1064 (2.38-2.34)
Ramachandran outliers	78287	1048 (2.38-2.34)
Sidechain outliers	78261	1049 (2.38-2.34)
RSRZ outliers	66119	3330 (2.40-2.32)

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  $\geq 3$ , 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	287	
1	B	287	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4637 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 Shikimate 5'-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2200	1417	362	412	9			
1	B	287	Total	C	N	O	S	0	0	0
			2200	1417	362	412	9			

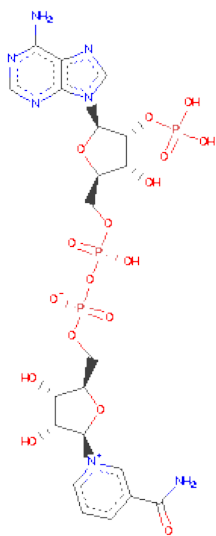
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP Q58484
A	2	PRO	-	CLONING ARTIFACT	UNP Q58484
A	3	LEU	-	CLONING ARTIFACT	UNP Q58484
A	4	GLY	-	CLONING ARTIFACT	UNP Q58484
A	5	SER	-	CLONING ARTIFACT	UNP Q58484
B	1	GLY	-	CLONING ARTIFACT	UNP Q58484
B	2	PRO	-	CLONING ARTIFACT	UNP Q58484
B	3	LEU	-	CLONING ARTIFACT	UNP Q58484
B	4	GLY	-	CLONING ARTIFACT	UNP Q58484
B	5	SER	-	CLONING ARTIFACT	UNP Q58484

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula: C<sub>21</sub>H<sub>28</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



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

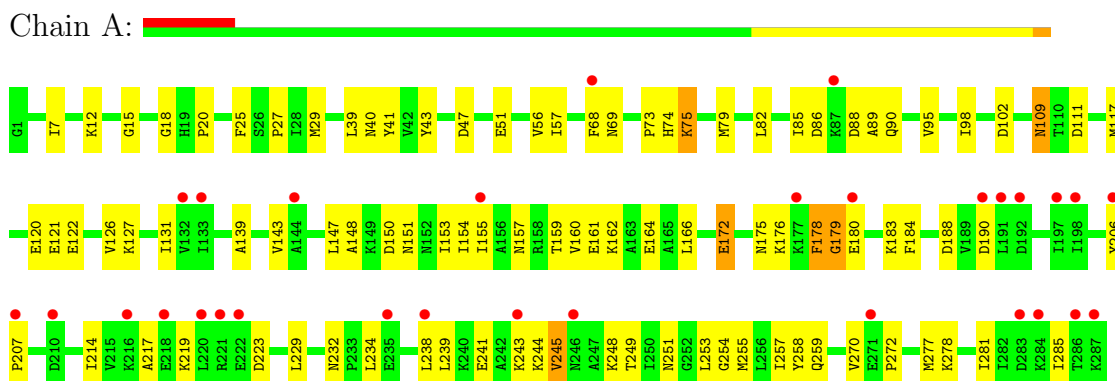
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	45	Total	O	0	0
			45	45		
4	B	94	Total	O	0	0
			94	94		

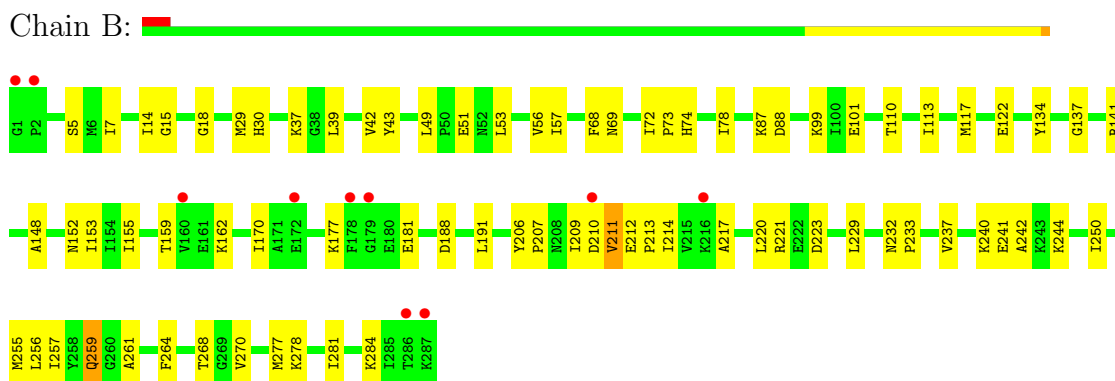
### 3 Residue-property plots

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: Shikimate 5'-dehydrogenase



- Molecule 1: Shikimate 5'-dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.65Å 75.62Å 118.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.96 – 2.35 21.96 – 2.35	Depositor EDS
% Data completeness (in resolution range)	97.9 (21.96-2.35) 99.1 (21.96-2.35)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.10 (at 2.36Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.224 , 0.260 0.230 , 0.269	Depositor DCC
$R_{free}$ test set	1398 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.4	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 27.5	EDS
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 53801 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4637	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.33	0/2232	0.60	1/3012 (0.0%)
1	B	0.37	0/2232	0.63	1/3012 (0.0%)
All	All	0.35	0/4464	0.62	2/6024 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	ASN	N-CA-C	-5.84	95.24	111.00
1	B	69	ASN	N-CA-C	-5.52	96.10	111.00

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	2200	0	2295	73	0
1	B	2200	0	2295	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	48	0	22	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	48	0	23	1	0
4	A	45	0	0	1	0
4	B	94	0	0	1	0
All	All	4637	0	4635	127	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 14.

All (127) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:268:THR:HG23	1:B:270:VAL:H	1.36	0.91
1:A:127:LYS:HG3	1:A:150:ASP:HB2	1.58	0.86
1:A:29:MET:HB2	1:A:281:ILE:HD11	1.64	0.79
1:A:25:PHE:HB3	1:A:281:ILE:HD12	1.64	0.79
1:B:29:MET:HE2	1:B:281:ILE:HD11	1.71	0.73
1:A:147:LEU:O	1:A:151:ASN:HB2	1.89	0.72
1:A:148:ALA:HB2	1:A:153:ILE:HD13	1.71	0.72
1:A:117:MET:HE3	1:A:258:TYR:HB3	1.74	0.70
1:B:257:ILE:HD11	1:B:278:LYS:HA	1.74	0.70
1:A:75:LYS:HB3	1:A:95:VAL:O	1.92	0.69
1:B:137:GLY:O	1:B:141:ARG:HG3	1.92	0.68
1:B:211:VAL:HG22	1:B:212:GLU:H	1.59	0.68
1:B:87:LYS:HG3	1:B:88:ASP:H	1.58	0.68
1:B:264:PHE:O	1:B:268:THR:HG22	1.95	0.67
1:A:12:LYS:HE2	1:A:12:LYS:HA	1.76	0.67
1:B:122:GLU:HG2	1:B:250:ILE:HG12	1.76	0.65
1:B:211:VAL:HG22	1:B:212:GLU:N	2.11	0.65
1:A:239:LEU:HD22	1:A:249:THR:HB	1.78	0.64
1:A:7:ILE:HD13	1:B:42:VAL:CG2	2.28	0.64
1:A:126:VAL:HG12	1:A:151:ASN:HD21	1.63	0.63
1:B:134:TYR:HB3	1:B:214:ILE:HD13	1.82	0.62
1:B:220:LEU:HD13	1:B:242:ALA:HA	1.80	0.62
1:A:51:GLU:H	1:A:51:GLU:CD	2.04	0.61
1:A:272:PRO:CG	1:A:277:MET:HE2	2.32	0.60
1:B:241:GLU:O	1:B:244:LYS:HG2	2.02	0.59
1:B:268:THR:HG23	1:B:270:VAL:N	2.14	0.59
1:A:229:LEU:HD22	1:A:255:MET:SD	2.43	0.58
1:B:237:VAL:HA	1:B:240:LYS:HE2	1.86	0.58
1:B:53:LEU:O	1:B:56:VAL:HG22	2.04	0.57
1:A:117:MET:CE	1:A:258:TYR:HB3	2.33	0.57
1:B:152:ASN:HD22	1:B:181:GLU:HA	1.70	0.57
1:A:206:TYR:CD1	1:A:207:PRO:HA	2.39	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:39:LEU:HD11	1:A:270:VAL:HG21	1.87	0.56
1:A:154:ILE:HG12	1:A:183:LYS:HB2	1.88	0.56
1:B:15:GLY:O	1:B:68:PHE:HA	2.07	0.55
1:A:117:MET:HE3	1:A:258:TYR:CD1	2.41	0.55
1:A:179:GLY:O	1:A:183:LYS:HE2	2.07	0.55
1:B:148:ALA:HB2	1:B:153:ILE:HD13	1.89	0.54
1:B:206:TYR:CE2	1:B:207:PRO:HG3	2.42	0.54
1:B:261:ALA:HB2	1:B:277:MET:HE3	1.89	0.54
1:A:122:GLU:HG3	1:A:248:LYS:NZ	2.21	0.54
1:A:79:MET:HE3	1:A:85:ILE:HD13	1.90	0.54
1:B:221:ARG:HD2	1:B:223:ASP:OD1	2.07	0.54
1:A:217:ALA:HB2	1:A:241:GLU:OE2	2.08	0.54
1:B:53:LEU:HD11	1:B:78:ILE:HD13	1.90	0.54
1:B:177:LYS:N	1:B:177:LYS:HD2	2.24	0.53
1:B:37:LYS:HB3	1:B:39:LEU:HG	1.92	0.52
1:A:79:MET:CE	1:A:85:ILE:HD13	2.41	0.51
1:A:126:VAL:HG12	1:A:151:ASN:ND2	2.26	0.51
1:B:134:TYR:HB3	1:B:214:ILE:CD1	2.40	0.51
1:B:30:HIS:HE1	4:B:292:HOH:O	1.94	0.50
1:B:191:LEU:O	1:B:221:ARG:NH2	2.44	0.50
1:B:7:ILE:HD12	1:B:7:ILE:N	2.28	0.49
1:B:177:LYS:HA	1:B:177:LYS:HE3	1.95	0.48
1:A:175:ASN:O	1:A:176:LYS:HD2	2.14	0.47
1:A:127:LYS:NZ	1:A:127:LYS:HB3	2.28	0.47
1:B:29:MET:HE1	1:B:256:LEU:HG	1.96	0.47
1:A:254:GLY:O	1:A:258:TYR:HD2	1.98	0.47
1:B:99:LYS:HE2	1:B:101:GLU:OE2	2.14	0.47
1:B:261:ALA:HB2	1:B:277:MET:CE	2.45	0.46
1:A:86:ASP:O	1:A:89:ALA:HB3	2.15	0.46
1:A:85:ILE:HG22	1:A:90:GLN:HG2	1.98	0.46
1:A:15:GLY:O	1:A:68:PHE:HA	2.14	0.46
1:A:109:ASN:HD21	1:A:111:ASP:HB2	1.81	0.45
1:A:131:ILE:O	1:A:153:ILE:HA	2.16	0.45
1:A:120:GLU:O	1:A:122:GLU:N	2.49	0.45
1:A:18:GLY:HA2	1:A:74:HIS:CD2	2.51	0.45
1:A:109:ASN:C	1:A:109:ASN:HD22	2.20	0.45
1:A:253:LEU:HG	1:A:257:ILE:HD12	1.99	0.45
1:A:117:MET:HE3	1:A:258:TYR:CB	2.45	0.44
1:A:184:PHE:CD1	1:A:184:PHE:C	2.90	0.44
1:A:234:LEU:HD11	1:A:285:ILE:HD13	1.99	0.44
1:A:190:ASP:HA	1:A:219:LYS:NZ	2.33	0.44
1:B:113:ILE:HG22	1:B:117:MET:CE	2.48	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:253:LEU:HG	1:A:257:ILE:CD1	2.47	0.44
1:B:211:VAL:CG2	1:B:212:GLU:H	2.28	0.44
1:B:209:ILE:HG21	1:B:233:PRO:HG3	2.00	0.44
1:B:155:ILE:HD11	1:B:170:ILE:HD12	2.00	0.44
1:B:18:GLY:HA2	1:B:74:HIS:CD2	2.52	0.44
1:B:122:GLU:HG2	1:B:250:ILE:CG1	2.47	0.44
1:A:39:LEU:HB3	1:A:41:TYR:HD2	1.81	0.44
1:A:164:GLU:HG3	1:A:184:PHE:CE2	2.53	0.44
1:B:159:THR:O	1:B:162:LYS:HB3	2.18	0.44
1:A:148:ALA:HB2	1:A:153:ILE:CD1	2.45	0.43
1:A:120:GLU:C	1:A:122:GLU:H	2.21	0.43
1:A:272:PRO:CG	1:A:277:MET:CE	2.96	0.43
1:B:217:ALA:HB2	1:B:241:GLU:HG3	1.98	0.43
1:B:229:LEU:HD22	1:B:255:MET:SD	2.58	0.43
1:A:139:ALA:O	1:A:143:VAL:HG23	2.18	0.43
1:B:213:PRO:HG3	1:B:237:VAL:CG2	2.48	0.43
1:B:49:LEU:HD13	1:B:51:GLU:OE2	2.18	0.43
1:A:178:PHE:O	1:A:180:GLU:N	2.51	0.43
1:A:214:ILE:HG22	1:A:238:LEU:HD13	2.00	0.43
1:B:110:THR:OG1	1:B:259:GLN:HG2	2.19	0.43
1:A:175:ASN:C	1:A:176:LYS:HD2	2.39	0.42
1:A:157:ASN:ND2	4:A:296:HOH:O	2.52	0.42
1:B:14:ILE:O	1:B:43:TYR:HA	2.19	0.42
1:B:214:ILE:O	1:B:214:ILE:HG12	2.20	0.42
1:B:113:ILE:HG22	1:B:117:MET:HE1	2.00	0.42
1:B:72:ILE:HD11	3:B:289:NAP:O1N	2.18	0.42
1:A:82:LEU:HD21	1:A:98:ILE:HD13	2.02	0.42
1:A:120:GLU:C	1:A:122:GLU:N	2.72	0.42
1:B:56:VAL:HG23	1:B:57:ILE:N	2.34	0.42
1:A:20:PRO:HA	1:A:47:ASP:OD1	2.20	0.42
1:A:244:LYS:O	1:A:245:VAL:HG13	2.19	0.42
1:B:211:VAL:CG2	1:B:212:GLU:N	2.79	0.41
1:B:217:ALA:N	1:B:241:GLU:OE1	2.45	0.41
1:A:160:VAL:HG23	1:A:161:GLU:N	2.36	0.41
1:A:122:GLU:HG3	1:A:248:LYS:HZ1	1.84	0.41
1:B:29:MET:CE	1:B:256:LEU:HG	2.50	0.41
1:A:27:PRO:HA	1:A:43:TYR:CD2	2.55	0.41
1:A:272:PRO:HG2	1:A:277:MET:CE	2.51	0.41
1:B:5:SER:O	1:B:7:ILE:HD12	2.20	0.41
1:A:251:ASN:HD21	1:A:253:LEU:HB2	1.86	0.41
1:A:253:LEU:HD21	1:A:278:LYS:HG3	2.01	0.41
1:A:234:LEU:N	1:A:234:LEU:HD12	2.36	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:7:ILE:CD1	1:B:42:VAL:CG2	2.98	0.41
1:B:213:PRO:HG3	1:B:237:VAL:HB	2.03	0.41
1:A:86:ASP:O	1:A:90:GLN:HG3	2.21	0.41
1:A:243:LYS:C	1:A:245:VAL:H	2.23	0.41
1:A:223:ASP:OD2	1:A:223:ASP:N	2.54	0.41
1:A:39:LEU:HB3	1:A:41:TYR:CD2	2.56	0.40
1:A:155:ILE:HD13	1:A:166:LEU:HD23	2.02	0.40
1:A:56:VAL:HG23	1:A:57:ILE:N	2.36	0.40
1:A:159:THR:HG23	3:A:288:NAP:O2X	2.22	0.40
1:A:172:GLU:C	1:A:172:GLU:OE1	2.60	0.40
1:A:85:ILE:CG2	1:A:90:GLN:HG2	2.52	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	285/287 (99%)	258 (90%)	23 (8%)	4 (1%)	16	15
1	B	285/287 (99%)	265 (93%)	19 (7%)	1 (0%)	43	52
All	All	570/574 (99%)	523 (92%)	42 (7%)	5 (1%)	25	27

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	178	PHE
1	A	179	GLY
1	A	121	GLU
1	A	245	VAL
1	B	211	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	233/233 (100%)	222 (95%)	11 (5%)	36	46
1	B	233/233 (100%)	227 (97%)	6 (3%)	59	75
All	All	466/466 (100%)	449 (96%)	17 (4%)	47	61

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	73	PRO
1	A	75	LYS
1	A	88	ASP
1	A	102	ASP
1	A	109	ASN
1	A	162	LYS
1	A	172	GLU
1	A	188	ASP
1	A	232	ASN
1	A	259	GLN
1	B	73	PRO
1	B	188	ASP
1	B	210	ASP
1	B	232	ASN
1	B	259	GLN
1	B	284	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	40	ASN
1	A	52	ASN
1	A	109	ASN
1	A	130	ASN
1	A	152	ASN
1	A	157	ASN

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Mol	Chain	Res	Type
1	A	232	ASN
1	A	279	ASN
1	B	30	HIS
1	B	52	ASN
1	B	152	ASN
1	B	175	ASN
1	B	199	ASN
1	B	208	ASN
1	B	232	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 ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z  > 2$
3	NAP	A	288	-	52,52,52	2.59	15 (28%)	80,80,80	4.47	23 (28%)
3	NAP	B	289	-	52,52,52	2.54	15 (28%)	80,80,80	4.47	24 (30%)

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
3	NAP	A	288	-	1/1/12/12	0/35/67/67	0/3/5/5
3	NAP	B	289	-	1/1/12/12	0/35/67/67	0/3/5/5

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	289	NAP	O7N-C7N	8.13	1.43	1.24
3	A	288	NAP	O7N-C7N	8.07	1.43	1.24
3	B	289	NAP	C2N-N1N	7.60	1.45	1.35
3	A	288	NAP	C2N-N1N	7.16	1.44	1.35
3	A	288	NAP	C5N-C4N	5.61	1.52	1.39
3	B	289	NAP	C5N-C4N	5.60	1.52	1.39
3	B	289	NAP	C6N-C5N	5.59	1.51	1.38
3	A	288	NAP	PA-O3	-5.49	1.49	1.59
3	A	288	NAP	PN-O3	-5.38	1.49	1.60
3	A	288	NAP	C6N-C5N	5.22	1.50	1.38
3	B	289	NAP	PA-O3	-4.57	1.51	1.59
3	B	289	NAP	PN-O3	-4.23	1.51	1.60
3	A	288	NAP	C6N-N1N	3.90	1.46	1.35
3	B	289	NAP	C6N-N1N	3.84	1.46	1.35
3	B	289	NAP	C4N-C3N	3.65	1.45	1.39
3	A	288	NAP	C4N-C3N	3.55	1.45	1.39
3	A	288	NAP	C1D-N1N	-3.47	1.37	1.48
3	B	289	NAP	C1D-N1N	-3.43	1.38	1.48
3	A	288	NAP	C2N-C3N	3.39	1.43	1.38
3	B	289	NAP	C2N-C3N	2.97	1.43	1.38
3	A	288	NAP	C1B-N9A	-2.86	1.39	1.48
3	B	289	NAP	C4A-N3A	2.69	1.39	1.35
3	B	289	NAP	C2A-N3A	2.64	1.37	1.32
3	A	288	NAP	C2A-N3A	2.47	1.37	1.32
3	A	288	NAP	C4A-N3A	2.47	1.39	1.35
3	A	288	NAP	PN-O5D	-2.43	1.52	1.60
3	B	289	NAP	C2A-N1A	2.36	1.38	1.33
3	B	289	NAP	C1B-N9A	-2.35	1.41	1.48
3	A	288	NAP	C2A-N1A	2.30	1.38	1.33
3	B	289	NAP	PN-O5D	-2.26	1.52	1.60

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	288	NAP	C2D-C1D-N1N	24.86	155.95	113.86
3	B	289	NAP	C2D-C1D-N1N	24.48	155.30	113.86
3	B	289	NAP	P2B-O2B-C2B	-17.75	84.58	121.96
3	A	288	NAP	P2B-O2B-C2B	-16.79	86.61	121.96
3	A	288	NAP	O4D-C1D-N1N	-11.42	96.27	107.95
3	A	288	NAP	N3A-C2A-N1A	-10.69	119.77	128.71
3	B	289	NAP	O4D-C1D-N1N	-10.53	97.18	107.95
3	B	289	NAP	N3A-C2A-N1A	-10.46	119.96	128.71
3	A	288	NAP	PN-O5D-C5D	8.18	148.51	120.24
3	B	289	NAP	PN-O5D-C5D	7.92	147.64	120.24
3	B	289	NAP	O5D-C5D-C4D	6.92	134.32	108.94
3	A	288	NAP	O5D-C5D-C4D	6.68	133.45	108.94
3	B	289	NAP	O3X-P2B-O2B	6.57	126.02	107.09
3	A	288	NAP	O3X-P2B-O2B	6.51	125.84	107.09
3	A	288	NAP	C6N-N1N-C1D	-6.51	103.21	119.33
3	B	289	NAP	C6N-N1N-C1D	-6.28	103.77	119.33
3	B	289	NAP	O4B-C1B-N9A	-6.01	102.85	108.44
3	A	288	NAP	C2N-N1N-C1D	6.00	134.20	119.33
3	B	289	NAP	C2N-N1N-C1D	5.76	133.60	119.33
3	B	289	NAP	N3A-C4A-N9A	5.57	135.48	125.43
3	A	288	NAP	O4B-C1B-N9A	-5.54	103.29	108.44
3	A	288	NAP	N3A-C4A-N9A	5.46	135.29	125.43
3	B	289	NAP	C2B-C1B-N9A	5.43	134.33	113.74
3	A	288	NAP	C2B-C1B-N9A	5.08	132.99	113.74
3	A	288	NAP	C3D-C2D-C1D	4.10	107.32	100.91
3	B	289	NAP	C3D-C2D-C1D	4.07	107.28	100.91
3	B	289	NAP	O3-PN-O5D	3.81	116.63	101.36
3	A	288	NAP	O3-PN-O5D	3.44	115.14	101.36
3	A	288	NAP	O2X-P2B-O2B	-3.20	97.87	107.09
3	B	289	NAP	C5A-C4A-N3A	-3.16	118.82	125.70
3	A	288	NAP	O2N-PN-O1N	-3.05	109.51	118.72
3	A	288	NAP	C5A-C4A-N3A	-2.98	119.22	125.70
3	B	289	NAP	O4B-C1B-C2B	-2.93	104.20	106.95
3	B	289	NAP	O2X-P2B-O2B	-2.85	98.89	107.09
3	B	289	NAP	O2N-PN-O1N	-2.78	110.34	118.72
3	A	288	NAP	C8A-N9A-C1B	-2.76	120.94	126.38
3	B	289	NAP	C1B-N9A-C4A	2.66	131.23	126.64
3	A	288	NAP	C2A-N3A-C4A	2.63	121.50	114.01
3	B	289	NAP	C2A-N3A-C4A	2.63	121.50	114.01
3	B	289	NAP	C3B-C2B-C1B	-2.47	97.94	102.73
3	B	289	NAP	C6N-C5N-C4N	-2.39	115.63	119.44
3	B	289	NAP	C4A-C5A-N7A	2.39	111.57	109.52
3	A	288	NAP	C6N-C5N-C4N	-2.38	115.65	119.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	288	NAP	C1B-N9A-C4A	2.35	130.69	126.64
3	B	289	NAP	C8A-N9A-C1B	-2.31	121.84	126.38
3	A	288	NAP	C4A-C5A-N7A	2.26	111.46	109.52
3	A	288	NAP	C3B-C2B-C1B	-2.17	98.52	102.73

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	288	NAP	C1D
3	B	289	NAP	C1D

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	287/287 (100%)	0.67	30 (10%) 7 8	23, 49, 81, 92	0
1	B	287/287 (100%)	0.20	10 (3%) 42 46	19, 34, 60, 70	0
All	All	574/574 (100%)	0.43	40 (6%) 17 18	19, 42, 77, 92	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	286	THR	5.9
1	A	287	LYS	5.3
1	A	246	ASN	4.8
1	A	243	LYS	4.6
1	A	210	ASP	4.6
1	B	287	LYS	4.1
1	A	133	ILE	3.9
1	A	283	ASP	3.4
1	A	177	LYS	3.2
1	A	206	TYR	3.1
1	A	144	ALA	3.1
1	A	192	ASP	3.1
1	A	198	ILE	2.8
1	A	87	LYS	2.8
1	B	210	ASP	2.8
1	A	238	LEU	2.8
1	B	1	GLY	2.7
1	B	172	GLU	2.7
1	A	191	LEU	2.6
1	A	207	PRO	2.6
1	A	235	GLU	2.5
1	A	197	ILE	2.5
1	A	132	VAL	2.5
1	B	179	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	220	LEU	2.4
1	A	221	ARG	2.4
1	B	178	PHE	2.4
1	A	284	LYS	2.3
1	A	218	GLU	2.3
1	A	222	GLU	2.3
1	A	190	ASP	2.3
1	B	2	PRO	2.3
1	A	180	GLU	2.1
1	A	68	PHE	2.1
1	A	271	GLU	2.1
1	B	286	THR	2.1
1	A	216	LYS	2.0
1	A	155	ILE	2.0
1	B	160	VAL	2.0
1	B	216	LYS	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 ⓘ

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, 95<sup>th</sup> 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(Å <sup>2</sup> )	Q<0.9
3	NAP	B	289	48/48	0.18	1.08	50,64,68,69	0
3	NAP	A	288	48/48	0.14	0.17	38,48,54,55	0
2	ZN	B	291	1/1	0.06	-1.97	36,36,36,36	0
2	ZN	A	290	1/1	0.05	-2.35	46,46,46,46	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.