



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:42 am GMT

PDB ID : 1KEV  
Title : STRUCTURE OF NADP-DEPENDENT ALCOHOL DEHYDROGENASE  
Authors : Korkhin, Y.; Frolov, F.  
Deposited on : 1996-10-21  
Resolution : 2.05 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](mailto: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.

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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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

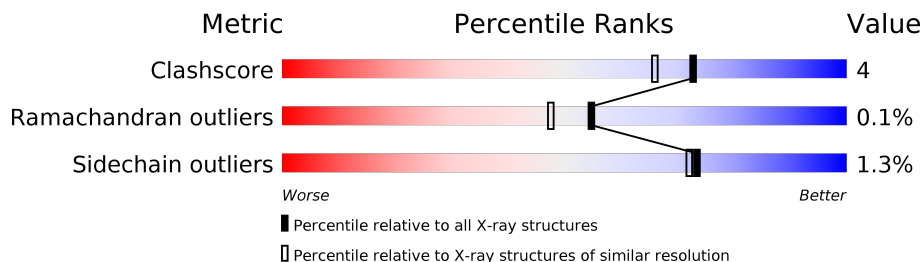
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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))
Clashscore	112137	1394 (2.04-2.04)
Ramachandran outliers	110173	1383 (2.04-2.04)
Sidechain outliers	110143	1383 (2.04-2.04)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	351	<div> <div style="width: 91%;"></div> <div style="width: 8%;"></div> <div style="width: 1%;"></div> </div> <div>91% 8% .</div>
1	B	351	<div> <div style="width: 94%;"></div> <div style="width: 6%;"></div> </div> <div>94% 6%</div>
1	C	351	<div> <div style="width: 89%;"></div> <div style="width: 11%;"></div> </div> <div>89% 11%</div>
1	D	351	<div> <div style="width: 91%;"></div> <div style="width: 9%;"></div> </div> <div>91% 9%</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
3	NDP	A	352	X	-	-	-
3	NDP	B	352	X	-	-	-
3	NDP	C	352	X	-	-	-
3	NDP	D	352	X	-	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11442 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 NADP-DEPENDENT ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	B	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	C	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			
1	D	351	Total	C	N	O	S	0	0	0
			2640	1675	460	482	23			

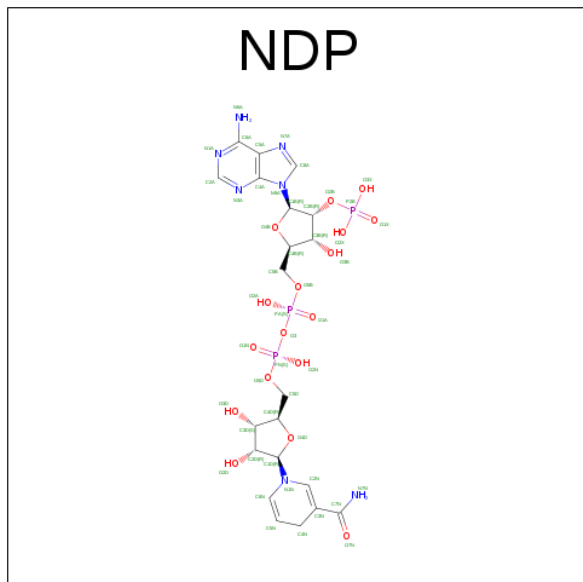
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	THR	SER	CONFLICT	UNP P25984
A	234	LYS	GLU	CONFLICT	UNP P25984
B	154	THR	SER	CONFLICT	UNP P25984
B	234	LYS	GLU	CONFLICT	UNP P25984
C	154	THR	SER	CONFLICT	UNP P25984
C	234	LYS	GLU	CONFLICT	UNP P25984
D	154	THR	SER	CONFLICT	UNP P25984
D	234	LYS	GLU	CONFLICT	UNP P25984

- 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		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula:  $C_{21}H_{30}N_7O_{17}P_3$ ).



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		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is water.

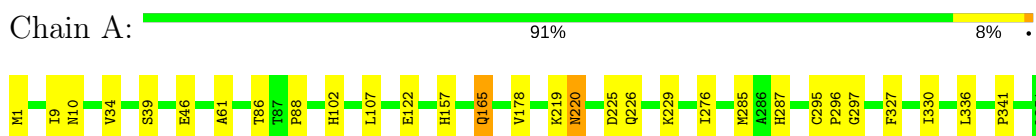
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	180	Total	O	0	0
			180	180		
4	B	177	Total	O	0	0
			177	177		
4	C	154	Total	O	0	0
			154	154		
4	D	175	Total	O	0	0
			175	175		

### 3 Residue-property plots [i](#)

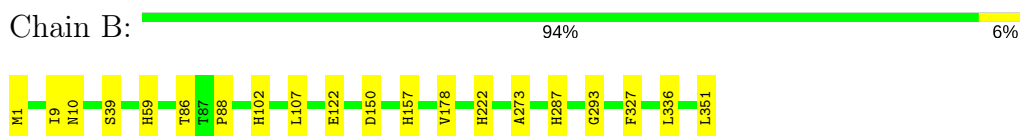
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

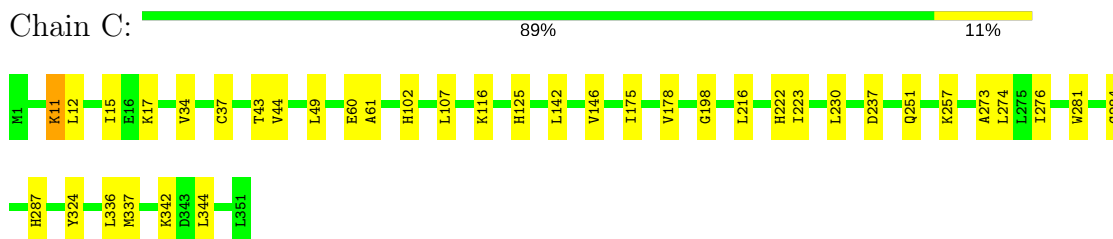
#### • Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



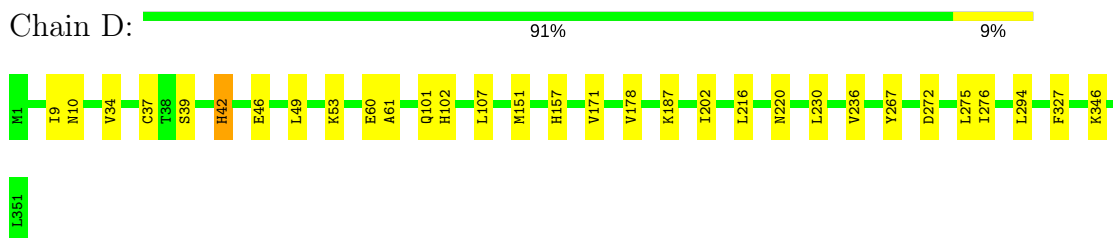
#### • Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



#### • Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



#### • Molecule 1: NADP-DEPENDENT ALCOHOL DEHYDROGENASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.45Å 151.42Å 127.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.05	Depositor
% Data completeness (in resolution range)	88.2 (50.00-2.05)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.205 , 0.259	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	11442	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

## 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, ZN

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	0.34	0/2689	0.60	0/3626
1	B	0.36	0/2689	0.60	0/3626
1	C	0.34	0/2689	0.60	0/3626
1	D	0.35	0/2689	0.60	0/3626
All	All	0.35	0/10756	0.60	0/14504

There are no bond length outliers.

There are no bond angle outliers.

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	2640	0	2681	28	0
1	B	2640	0	2681	22	0
1	C	2640	0	2681	31	0
1	D	2640	0	2681	24	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	48	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	48	0	24	3	0
3	C	48	0	25	1	0
3	D	48	0	24	1	0
4	A	180	0	0	1	0
4	B	177	0	0	0	0
4	C	154	0	0	2	0
4	D	175	0	0	2	0
All	All	11442	0	10821	84	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 4.

All (84) 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:178:VAL:HG21	3:A:352:NDP:H6N	1.43	1.01
1:C:178:VAL:HG21	3:C:352:NDP:H6N	1.59	0.84
1:D:178:VAL:HG21	3:D:352:NDP:H6N	1.59	0.83
1:D:102:HIS:HD2	1:D:107:LEU:H	1.25	0.82
1:A:102:HIS:HD2	1:A:107:LEU:H	1.26	0.81
1:C:102:HIS:HD2	1:C:107:LEU:H	1.36	0.73
1:B:178:VAL:HG21	3:B:352:NDP:H6N	1.70	0.73
1:B:102:HIS:HD2	1:B:107:LEU:H	1.39	0.70
1:D:216:LEU:HD21	1:D:230:LEU:HD11	1.74	0.68
1:C:11:LYS:NZ	1:C:11:LYS:HA	2.09	0.67
1:C:11:LYS:HZ2	1:C:12:LEU:H	1.46	0.64
1:D:230:LEU:HD13	4:D:498:HOH:O	1.99	0.63
1:C:287:HIS:HE1	1:D:157:HIS:HE1	1.45	0.62
1:B:327:PHE:HA	1:B:351:LEU:HD22	1.83	0.60
1:A:102:HIS:HE1	1:B:287:HIS:HD2	1.47	0.60
1:C:11:LYS:HZ3	1:C:11:LYS:HA	1.64	0.59
1:A:157:HIS:HE1	1:B:287:HIS:HE1	1.48	0.58
1:D:9:ILE:HG22	1:D:10:ASN:HD22	1.68	0.58
1:D:102:HIS:CD2	1:D:107:LEU:H	2.15	0.58
1:A:102:HIS:CE1	1:B:287:HIS:HD2	2.21	0.57
1:D:151:MET:SD	1:D:346:LYS:HE2	2.46	0.55
1:C:34:VAL:HG12	1:C:61:ALA:HB2	1.87	0.55
1:B:9:ILE:HG22	1:B:10:ASN:ND2	2.21	0.55
1:D:53:LYS:HB2	1:D:53:LYS:NZ	2.22	0.55
1:A:9:ILE:HG22	1:A:10:ASN:ND2	2.21	0.54
1:B:1:MET:HG3	1:B:122:GLU:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:CYS:HB2	1:C:60:GLU:OE2	2.07	0.54
1:A:102:HIS:HE1	1:B:287:HIS:CD2	2.25	0.54
1:D:187:LYS:HD3	4:D:528:HOH:O	2.08	0.53
1:A:178:VAL:HG21	3:A:352:NDP:C6N	2.29	0.52
1:A:34:VAL:HG12	1:A:61:ALA:HB2	1.91	0.51
1:A:86:THR:HG22	1:A:88:PRO:HD3	1.91	0.51
1:A:287:HIS:CE1	1:B:157:HIS:HE1	2.28	0.51
1:C:337:MET:HE2	1:C:344:LEU:HD21	1.91	0.50
1:A:1:MET:HG3	1:A:122:GLU:HB2	1.94	0.50
1:A:287:HIS:HE1	1:B:157:HIS:HE1	1.58	0.50
1:C:287:HIS:CE1	1:D:157:HIS:HE1	2.28	0.50
1:D:42:HIS:O	1:D:46:GLU:HB2	2.11	0.50
1:C:216:LEU:HD21	1:C:230:LEU:HD11	1.94	0.50
1:C:287:HIS:HD2	1:D:102:HIS:HE1	1.61	0.49
1:B:150:ASP:OD1	3:B:352:NDP:H41N	2.12	0.48
1:C:273:ALA:HB1	1:D:276:ILE:O	2.13	0.48
1:C:125:HIS:CE1	4:C:445:HOH:O	2.65	0.48
1:A:86:THR:HB	1:A:297:GLY:HA3	1.95	0.48
1:A:220:ASN:HB3	1:A:226:GLN:CD	2.35	0.47
1:C:43:THR:HG22	1:C:49:LEU:HB2	1.95	0.47
1:D:101:GLN:HB3	1:D:294:LEU:HD23	1.97	0.47
1:C:287:HIS:CD2	1:D:102:HIS:HE1	2.33	0.47
1:A:102:HIS:CD2	1:A:107:LEU:H	2.17	0.47
1:A:157:HIS:HE1	1:B:287:HIS:CE1	2.30	0.46
1:D:102:HIS:HD2	1:D:107:LEU:N	2.05	0.46
1:D:171:VAL:HG23	1:D:236:VAL:HG11	1.97	0.46
1:A:287:HIS:HD2	1:B:102:HIS:HE1	1.64	0.46
1:C:337:MET:CE	1:C:344:LEU:HD21	2.46	0.46
1:C:116:LYS:NZ	1:C:125:HIS:HD2	2.14	0.46
1:C:324:TYR:OH	1:C:336:LEU:HD11	2.15	0.46
1:A:285:MET:HA	1:B:293:GLY:HA2	1.97	0.45
1:B:178:VAL:HG21	3:B:352:NDP:C6N	2.45	0.45
1:C:273:ALA:HB3	1:D:275:LEU:HD11	1.97	0.45
1:A:287:HIS:CD2	1:B:102:HIS:HE1	2.33	0.45
1:C:287:HIS:HD2	1:D:102:HIS:CE1	2.35	0.45
1:D:9:ILE:O	1:D:10:ASN:HB2	2.18	0.44
1:C:222:HIS:HB3	4:C:478:HOH:O	2.17	0.44
1:D:37:CYS:HB2	1:D:60:GLU:OE2	2.17	0.44
1:B:39:SER:OG	1:B:59:HIS:HE1	2.00	0.44
1:B:86:THR:HG22	1:B:88:PRO:HD3	1.99	0.44
1:A:287:HIS:HE1	1:B:157:HIS:CE1	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:GLN:NE2	4:A:437:HOH:O	2.50	0.43
1:C:15:ILE:HD12	1:C:17:LYS:HD3	1.99	0.43
1:A:327:PHE:O	1:A:330:ILE:HG13	2.19	0.43
1:A:276:ILE:O	1:B:273:ALA:HB1	2.18	0.43
1:D:34:VAL:HG12	1:D:61:ALA:HB2	2.00	0.43
1:A:336:LEU:HD22	1:A:341:PRO:HG2	2.00	0.42
1:A:287:HIS:HD2	1:B:102:HIS:CE1	2.38	0.42
1:C:281:TRP:CE3	1:C:284:GLY:HA2	2.54	0.42
1:C:237:ASP:OD1	1:C:257:LYS:HE3	2.20	0.41
1:C:276:ILE:HD11	1:D:276:ILE:HD11	2.02	0.41
1:A:225:ASP:HB3	1:A:229:LYS:NZ	2.35	0.41
1:A:295:CYS:HA	1:A:296:PRO:HD3	1.88	0.41
1:C:142:LEU:O	1:C:146:VAL:HG23	2.20	0.41
1:C:223:ILE:HD11	1:C:251:GLN:OE1	2.20	0.41
1:C:12:LEU:HD21	1:C:44:VAL:HG11	2.03	0.41
1:C:274:LEU:N	1:C:274:LEU:HD12	2.36	0.41
1:C:175:ILE:HG12	1:C:198:GLY:HA3	2.02	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	349/351 (99%)	333 (95%)	16 (5%)	0	100	100
1	B	349/351 (99%)	336 (96%)	13 (4%)	0	100	100
1	C	349/351 (99%)	334 (96%)	14 (4%)	1 (0%)	44	35
1	D	349/351 (99%)	334 (96%)	14 (4%)	1 (0%)	44	35
All	All	1396/1404 (99%)	1337 (96%)	57 (4%)	2 (0%)	55	47

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	342	LYS
1	D	49	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	279/279 (100%)	274 (98%)	5 (2%)	64	60
1	B	279/279 (100%)	277 (99%)	2 (1%)	87	87
1	C	279/279 (100%)	278 (100%)	1 (0%)	93	93
1	D	279/279 (100%)	272 (98%)	7 (2%)	53	46
All	All	1116/1116 (100%)	1101 (99%)	15 (1%)	73	72

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	SER
1	A	46	GLU
1	A	165	GLN
1	A	219	LYS
1	A	220	ASN
1	B	222	HIS
1	B	336	LEU
1	C	11	LYS
1	D	39	SER
1	D	42	HIS
1	D	202	ILE
1	D	220	ASN
1	D	267	TYR
1	D	272	ASP
1	D	327	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	102	HIS
1	A	125	HIS
1	A	157	HIS
1	A	165	GLN
1	A	268	HIS
1	A	287	HIS
1	A	322	HIS
1	B	10	ASN
1	B	102	HIS
1	B	157	HIS
1	B	220	ASN
1	B	232	ASN
1	B	287	HIS
1	C	102	HIS
1	C	125	HIS
1	C	220	ASN
1	C	287	HIS
1	D	10	ASN
1	D	102	HIS
1	D	125	HIS
1	D	157	HIS
1	D	220	ASN
1	D	287	HIS
1	D	329	HIS

### 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 ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	NDP	A	352	-	43,52,52	1.26	4 (9%)	49,80,80	2.63	20 (40%)
3	NDP	B	352	-	43,52,52	1.35	7 (16%)	49,80,80	2.38	17 (34%)
3	NDP	C	352	-	43,52,52	1.31	4 (9%)	49,80,80	2.25	18 (36%)
3	NDP	D	352	-	43,52,52	1.32	4 (9%)	49,80,80	2.26	14 (28%)

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	NDP	A	352	-	2/2/14/17	1/30/77/77	0/5/5/5
3	NDP	B	352	-	2/2/14/17	0/30/77/77	0/5/5/5
3	NDP	C	352	-	1/1/14/17	1/30/77/77	0/5/5/5
3	NDP	D	352	-	2/2/14/17	0/30/77/77	0/5/5/5

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	352	NDP	C4N-C5N	-4.23	1.39	1.49
3	A	352	NDP	C4N-C5N	-4.21	1.39	1.49
3	C	352	NDP	C4N-C5N	-4.09	1.40	1.49
3	B	352	NDP	C4N-C5N	-3.94	1.40	1.49
3	C	352	NDP	PA-O5B	2.00	1.67	1.59
3	A	352	NDP	O2D-C2D	2.08	1.47	1.43
3	B	352	NDP	C6N-C5N	2.12	1.37	1.33
3	A	352	NDP	C6N-C5N	2.20	1.37	1.33
3	B	352	NDP	O4B-C1B	2.23	1.44	1.41
3	D	352	NDP	C6N-C5N	2.28	1.37	1.33
3	A	352	NDP	O4B-C1B	2.28	1.44	1.41
3	B	352	NDP	O2D-C2D	2.29	1.48	1.43
3	C	352	NDP	C6N-C5N	2.29	1.37	1.33
3	B	352	NDP	C2N-C3N	2.30	1.41	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	352	NDP	O4B-C1B	2.30	1.44	1.41
3	C	352	NDP	O4B-C1B	2.36	1.44	1.41
3	B	352	NDP	C1D-N1N	2.39	1.53	1.46
3	D	352	NDP	O2D-C2D	2.49	1.48	1.43
3	B	352	NDP	C5D-C4D	2.78	1.60	1.51

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	352	NDP	N3A-C2A-N1A	-8.03	121.87	128.86
3	D	352	NDP	N3A-C2A-N1A	-7.65	122.19	128.86
3	A	352	NDP	N3A-C2A-N1A	-7.56	122.28	128.86
3	B	352	NDP	N3A-C2A-N1A	-7.27	122.52	128.86
3	B	352	NDP	C4B-O4B-C1B	-5.53	103.89	109.77
3	D	352	NDP	C1D-N1N-C6N	-4.69	110.58	120.77
3	A	352	NDP	C5D-C4D-C3D	-4.62	97.69	115.29
3	B	352	NDP	C1D-N1N-C6N	-4.28	111.47	120.77
3	C	352	NDP	C1D-N1N-C6N	-4.19	111.67	120.77
3	C	352	NDP	C4B-O4B-C1B	-3.79	105.73	109.77
3	A	352	NDP	C1D-N1N-C6N	-3.73	112.67	120.77
3	D	352	NDP	C4B-O4B-C1B	-3.71	105.82	109.77
3	A	352	NDP	O5D-PN-O1N	-3.31	95.89	109.25
3	B	352	NDP	O3D-C3D-C2D	-3.20	101.58	111.83
3	A	352	NDP	C4B-O4B-C1B	-3.09	106.47	109.77
3	A	352	NDP	C2D-C3D-C4D	-3.09	96.60	102.62
3	B	352	NDP	C3N-C2N-N1N	-3.07	118.63	123.08
3	D	352	NDP	O3D-C3D-C2D	-2.99	102.26	111.83
3	A	352	NDP	C3N-C2N-N1N	-2.92	118.84	123.08
3	A	352	NDP	C4D-O4D-C1D	-2.89	103.02	109.47
3	C	352	NDP	O5D-PN-O1N	-2.51	99.14	109.25
3	C	352	NDP	O2B-P2B-O1X	-2.40	99.86	109.26
3	D	352	NDP	C3N-C2N-N1N	-2.35	119.67	123.08
3	C	352	NDP	O2D-C2D-C1D	-2.26	102.32	109.96
3	C	352	NDP	C3N-C2N-N1N	-2.20	119.89	123.08
3	A	352	NDP	O3B-C3B-C2B	-2.09	105.24	111.18
3	B	352	NDP	C5D-C4D-C3D	-2.09	107.34	115.29
3	C	352	NDP	C5D-C4D-C3D	-2.02	107.59	115.29
3	A	352	NDP	O3X-P2B-O2X	2.01	115.71	107.61
3	B	352	NDP	O3X-P2B-O2X	2.03	115.78	107.61
3	A	352	NDP	O4B-C4B-C3B	2.03	109.21	105.17
3	A	352	NDP	O2N-PN-O1N	2.03	122.81	112.28
3	B	352	NDP	O4B-C1B-C2B	2.06	110.20	106.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	352	NDP	O5D-C5D-C4D	2.06	116.31	109.00
3	C	352	NDP	O5D-C5D-C4D	2.07	116.34	109.00
3	C	352	NDP	C1D-N1N-C2N	2.08	124.61	121.09
3	C	352	NDP	O4D-C4D-C3D	2.08	109.30	105.17
3	C	352	NDP	O4B-C4B-C3B	2.15	109.44	105.17
3	A	352	NDP	O5B-C5B-C4B	2.18	116.71	109.00
3	D	352	NDP	O4D-C4D-C5D	2.35	117.34	109.40
3	B	352	NDP	O4D-C4D-C5D	2.36	117.38	109.40
3	D	352	NDP	O2D-C2D-C3D	2.39	119.47	111.83
3	B	352	NDP	N6A-C6A-N1A	2.44	123.61	118.77
3	D	352	NDP	O4B-C4B-C3B	2.58	110.30	105.17
3	A	352	NDP	N6A-C6A-N1A	2.63	123.98	118.77
3	C	352	NDP	C3B-C2B-C1B	2.68	107.99	102.75
3	B	352	NDP	C1D-N1N-C2N	2.88	125.98	121.09
3	B	352	NDP	O2D-C2D-C3D	2.93	121.21	111.83
3	D	352	NDP	C1D-N1N-C2N	2.98	126.15	121.09
3	D	352	NDP	C4A-C5A-N7A	3.07	112.37	109.41
3	C	352	NDP	O2D-C2D-C3D	3.15	121.92	111.83
3	D	352	NDP	O5D-C5D-C4D	3.27	120.58	109.00
3	C	352	NDP	C4A-C5A-N7A	3.28	112.58	109.41
3	D	352	NDP	O2B-C2B-C3B	3.77	125.58	111.63
3	C	352	NDP	O4D-C1D-N1N	3.77	115.66	108.07
3	B	352	NDP	O2B-C2B-C3B	3.77	125.59	111.63
3	C	352	NDP	O2B-C2B-C1B	3.81	124.26	110.06
3	A	352	NDP	O4D-C1D-N1N	3.85	115.83	108.07
3	A	352	NDP	C4A-C5A-N7A	3.98	113.25	109.41
3	B	352	NDP	O2B-C2B-C1B	4.05	125.15	110.06
3	A	352	NDP	O2B-C2B-C1B	4.05	125.15	110.06
3	B	352	NDP	C4A-C5A-N7A	4.16	113.43	109.41
3	C	352	NDP	O2B-C2B-C3B	4.26	127.40	111.63
3	A	352	NDP	O2B-C2B-C3B	4.39	127.90	111.63
3	D	352	NDP	O4D-C1D-N1N	4.46	117.04	108.07
3	D	352	NDP	O2B-C2B-C1B	4.51	126.86	110.06
3	A	352	NDP	O4D-C4D-C3D	4.88	114.87	105.17
3	B	352	NDP	O4D-C1D-N1N	5.22	118.57	108.07
3	A	352	NDP	O5D-C5D-C4D	6.38	131.64	109.00

All (7) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	C	352	NDP	C2B
3	D	352	NDP	C2B

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Mol	Chain	Res	Type	Atom
3	D	352	NDP	C1B
3	B	352	NDP	C2B
3	B	352	NDP	C1B
3	A	352	NDP	C2B
3	A	352	NDP	C1B

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	352	NDP	P2B-O2B-C2B-C3B
3	C	352	NDP	P2B-O2B-C2B-C3B

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	352	NDP	2	0
3	B	352	NDP	3	0
3	C	352	NDP	1	0
3	D	352	NDP	1	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

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.