



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:26 am GMT

PDB ID : 1OG6  
Title : YDHF, AN ALDO-KETO REDUCTASE FROM E.COLI COMPLEXED WITH NADPH  
Authors : Jeudy, S.; Abergel, C.; Claverie, J.M.  
Deposited on : 2003-04-24  
Resolution : 2.80 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

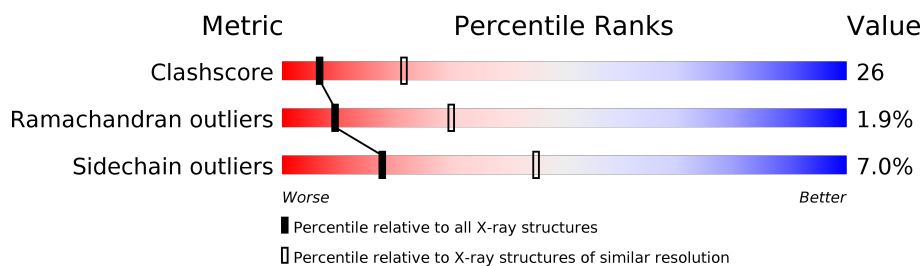
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)

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	298	
1	B	298	
1	C	298	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7423 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.

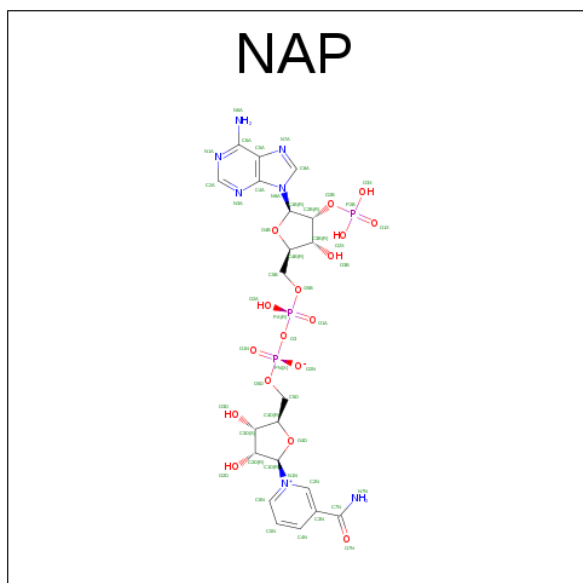
- Molecule 1 is a protein called HYPOTHETICAL OXIDOREDUCTASE YDHF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	1
			2362	1500	422	430	10			
1	B	294	Total	C	N	O	S	0	0	0
			2349	1493	418	428	10			
1	C	298	Total	C	N	O	S	0	0	0
			2369	1505	422	432	10			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	LEU	MET	CONFLICT	UNP P76187
B	1	LEU	MET	CONFLICT	UNP P76187
C	1	LEU	MET	CONFLICT	UNP P76187

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (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
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		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is water.

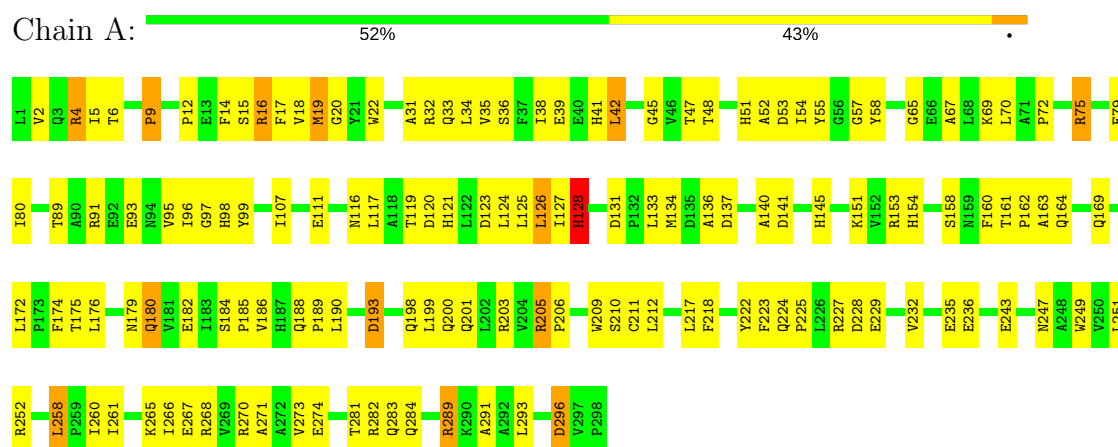
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total	O	0	0
			70	70		
3	B	76	Total	O	0	0
			76	76		
3	C	53	Total	O	0	0
			53	53		

### 3 Residue-property plots

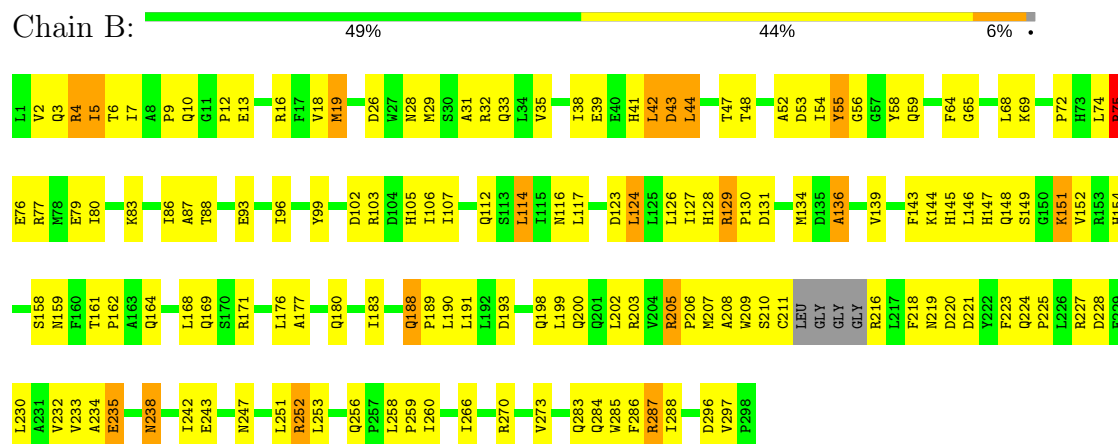
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

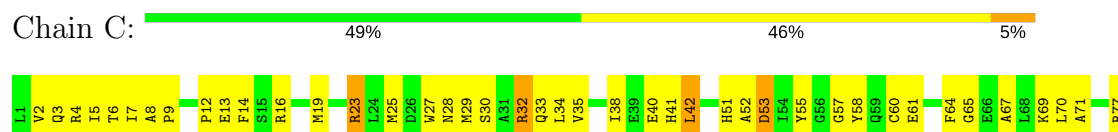
#### • Molecule 1: HYPOTHETICAL OXIDOREDUCTASE YDHF



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## 4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	174.53Å 174.53Å 97.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.91 – 2.80	Depositor
% Data completeness (in resolution range)	99.7 (29.91-2.80)	Depositor
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.207 , 0.238	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7423	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.36	0/2414	0.65	0/3278
1	B	0.37	0/2401	0.64	0/3259
1	C	0.36	0/2422	0.65	0/3288
All	All	0.36	0/7237	0.65	0/9825

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

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	2362	0	2355	123	0
1	B	2349	0	2341	124	0
1	C	2369	0	2362	146	0
2	A	48	0	25	8	0
2	B	48	0	25	7	0
2	C	48	0	25	5	0
3	A	70	0	0	4	0
3	B	76	0	0	6	0
3	C	53	0	0	13	0
All	All	7423	0	7133	375	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 26.

All (375) 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:C:161:THR:HB	1:C:164:GLN:HG3	1.40	1.02
1:A:95:VAL:H	1:C:283:GLN:NE2	1.60	1.00
1:A:210:SER:H	2:A:1298:NAP:H51N	1.28	0.96
1:B:75:ARG:HD3	1:B:75:ARG:O	1.67	0.94
1:C:88:THR:HG22	1:C:90:ALA:H	1.33	0.92
1:A:127:ILE:HD11	1:A:160:PHE:HZ	1.35	0.89
1:C:210:SER:H	2:C:1299:NAP:H51N	1.38	0.89
1:B:162:PRO:HA	1:B:198:GLN:HE22	1.37	0.87
1:A:2:VAL:HG12	1:A:45:GLY:HA3	1.56	0.85
1:B:129:ARG:HG3	1:B:130:PRO:HD2	1.60	0.84
1:C:169:GLN:HE22	1:C:176:LEU:H	1.23	0.83
1:A:95:VAL:H	1:C:283:GLN:HE22	1.22	0.83
1:C:250:VAL:HB	3:C:2038:HOH:O	1.76	0.83
1:C:38:ILE:O	1:C:42:LEU:HG	1.77	0.83
1:B:200:GLN:HE22	1:C:136:ALA:H	1.26	0.82
1:C:180:GLN:NE2	2:C:1299:NAP:H2N	1.94	0.81
1:C:112:GLN:HG3	1:C:116:ASN:HD21	1.45	0.81
1:A:18:VAL:CG2	1:A:260:ILE:HG12	2.10	0.80
1:B:149:SER:HB2	1:B:151:LYS:HG2	1.63	0.80
1:B:31:ALA:O	1:B:35:VAL:HG23	1.82	0.79
1:C:247:ASN:C	3:C:2038:HOH:O	2.21	0.78
1:B:54:ILE:HD11	3:B:2037:HOH:O	1.84	0.77
1:A:4:ARG:NH2	1:A:47:THR:OG1	2.17	0.77
1:A:162:PRO:HA	1:A:198:GLN:HE22	1.50	0.76
1:B:65:GLY:O	1:B:69:LYS:HG3	1.86	0.75
1:A:127:ILE:HD11	1:A:160:PHE:CZ	2.21	0.75
1:A:270:ARG:O	1:A:273:VAL:HG12	1.86	0.75
1:A:65:GLY:O	1:A:69:LYS:HG3	1.85	0.75
1:A:175:THR:HA	3:A:2036:HOH:O	1.86	0.74
1:B:284:GLN:O	1:B:288:ILE:HG12	1.87	0.74
1:C:218:PHE:HE2	1:C:243:GLU:HG2	1.53	0.74
1:B:38:ILE:O	1:B:42:LEU:HG	1.87	0.74
1:C:226:LEU:HD13	1:C:292:ALA:HB2	1.70	0.73
1:B:252:ARG:O	1:B:252:ARG:HD3	1.89	0.73
1:A:161:THR:HG22	1:A:163:ALA:H	1.52	0.72
1:C:144:LYS:HE2	1:C:173:PRO:HD3	1.69	0.71
1:A:121:HIS:CD2	1:A:153:ARG:HE	2.08	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:199:LEU:HD13	1:C:206:PRO:HB3	1.71	0.71
1:B:210:SER:H	2:B:1299:NAP:H51N	1.56	0.71
1:B:216:ARG:HA	1:B:219:ASN:HD22	1.55	0.71
1:C:103:ARG:HD3	1:C:104:ASP:N	2.06	0.71
1:B:224:GLN:HB3	1:B:225:PRO:HD3	1.72	0.71
1:C:121:HIS:CD2	1:C:153:ARG:HE	2.09	0.70
1:C:88:THR:HG22	1:C:90:ALA:N	2.06	0.70
1:B:162:PRO:HA	1:B:198:GLN:NE2	2.05	0.70
1:A:247:ASN:O	1:A:251:LEU:HD22	1.92	0.70
1:C:250:VAL:N	3:C:2038:HOH:O	2.24	0.70
1:C:7:ILE:HD13	1:C:14:PHE:CE2	2.26	0.70
1:A:169:GLN:HE22	1:A:176:LEU:H	1.38	0.70
1:C:111:GLU:HA	1:C:114:LEU:HD23	1.72	0.69
1:C:23:ARG:HH21	1:C:23:ARG:HG2	1.54	0.69
1:B:209:TRP:HB2	2:B:1299:NAP:O4D	1.93	0.69
1:B:180:GLN:NE2	2:B:1299:NAP:H2N	2.06	0.69
1:A:161:THR:HB	1:A:164:GLN:HG3	1.76	0.68
1:C:65:GLY:O	1:C:69:LYS:HG3	1.94	0.68
1:A:119:THR:HG22	1:A:120:ASP:N	2.08	0.68
1:A:121:HIS:HD2	1:A:153:ARG:HE	1.39	0.68
1:B:220:ASP:HB3	1:B:223:PHE:HD1	1.58	0.68
1:A:75:ARG:HH11	1:A:80:ILE:HD12	1.58	0.68
1:C:216:ARG:O	1:C:220:ASP:HB2	1.94	0.67
1:A:99:TYR:HB2	1:A:131:ASP:OD1	1.94	0.67
1:B:260:ILE:HG21	2:B:1299:NAP:O3D	1.95	0.67
1:A:18:VAL:HG21	1:A:260:ILE:HG12	1.76	0.66
1:A:119:THR:HG22	1:A:121:HIS:H	1.61	0.65
1:C:34:LEU:HD23	1:C:67:ALA:HB2	1.79	0.65
1:B:114:LEU:HD21	1:B:151:LYS:HB2	1.79	0.65
1:A:119:THR:HG22	1:A:120:ASP:H	1.62	0.64
1:B:251:LEU:HD13	1:B:259:PRO:HG2	1.80	0.64
1:C:144:LYS:HE2	1:C:173:PRO:CD	2.27	0.64
1:C:91:ARG:HB3	1:C:93:GLU:OE1	1.99	0.63
1:C:180:GLN:HE21	2:C:1299:NAP:H2N	1.63	0.63
1:C:189:PRO:O	1:C:192:LEU:HD23	1.99	0.63
1:A:127:ILE:O	1:A:127:ILE:HG13	1.97	0.63
1:A:39:GLU:O	1:A:42:LEU:HD12	1.98	0.63
1:B:189:PRO:HG2	1:B:190:LEU:HD22	1.80	0.63
1:A:128:HIS:O	1:A:158:SER:HB3	1.99	0.62
1:B:297:VAL:HG22	3:C:2014:HOH:O	1.99	0.62
1:C:16:ARG:HH21	1:C:16:ARG:HG3	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:HIS:HD2	1:C:153:ARG:HE	1.47	0.61
1:A:200:GLN:HE22	1:B:136:ALA:H	1.48	0.61
1:C:218:PHE:CE2	1:C:243:GLU:HG2	2.35	0.61
1:B:169:GLN:HE22	1:B:176:LEU:H	1.49	0.61
1:B:158:SER:O	1:B:159:ASN:HB2	2.01	0.61
1:C:53:ASP:HB2	1:C:84:CYS:HA	1.83	0.61
1:A:2:VAL:HG12	1:A:45:GLY:CA	2.31	0.61
1:C:181:VAL:HB	1:C:190:LEU:CD1	2.31	0.61
1:C:283:GLN:O	1:C:287:ARG:HB2	2.01	0.60
1:B:180:GLN:HE21	2:B:1299:NAP:H2N	1.64	0.60
1:C:212:LEU:HD13	1:C:218:PHE:HZ	1.67	0.60
1:B:233:VAL:HG21	1:B:288:ILE:HD11	1.85	0.59
1:B:3:GLN:HG2	1:B:4:ARG:N	2.16	0.59
1:C:281:THR:OG1	1:C:284:GLN:HB2	2.03	0.59
1:B:235:GLU:HG2	3:B:2062:HOH:O	2.02	0.59
1:C:199:LEU:CD1	1:C:206:PRO:HB3	2.33	0.58
1:B:183:ILE:HD12	1:B:208:ALA:HB2	1.86	0.58
1:A:210:SER:N	2:A:1298:NAP:H51N	2.10	0.58
1:A:14:PHE:CD2	1:A:48:THR:HG21	2.38	0.58
1:A:31:ALA:O	1:A:35:VAL:HG23	2.02	0.58
1:A:265:LYS:HB3	1:A:267:GLU:OE1	2.03	0.58
1:A:51:HIS:O	1:A:52:ALA:HB2	2.04	0.58
1:C:140:ALA:O	1:C:144:LYS:HG2	2.04	0.58
1:A:12:PRO:HB3	1:A:154:HIS:CD2	2.38	0.58
1:A:140:ALA:HA	1:A:172:LEU:HD21	1.85	0.58
1:A:79:GLU:HA	1:A:123:ASP:OD2	2.03	0.58
1:B:4:ARG:NH2	1:B:47:THR:OG1	2.36	0.57
1:A:217:LEU:HD23	1:A:217:LEU:O	2.05	0.57
1:A:229:GLU:HG3	1:A:291:ALA:HB2	1.86	0.57
1:A:184:SER:HB2	1:A:211:CYS:O	2.04	0.57
1:B:52:ALA:HB3	1:B:55:TYR:CD2	2.40	0.57
1:A:69:LYS:O	1:A:72:PRO:HD3	2.05	0.56
1:B:260:ILE:HB	2:B:1299:NAP:H4D	1.87	0.56
1:B:29:MET:HA	1:B:33:GLN:NE2	2.21	0.56
1:B:80:ILE:HG21	1:B:117:LEU:HD13	1.87	0.56
1:A:34:LEU:HD23	1:A:67:ALA:HB2	1.87	0.56
1:C:110:ALA:O	1:C:114:LEU:HD22	2.06	0.56
1:C:162:PRO:HA	1:C:198:GLN:HE22	1.70	0.56
1:A:91:ARG:HD3	1:A:93:GLU:OE1	2.06	0.56
1:C:103:ARG:C	1:C:103:ARG:HD3	2.27	0.55
1:A:35:VAL:O	1:A:39:GLU:HB2	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:64:PHE:HD2	1:C:117:LEU:HD22	1.71	0.55
1:A:35:VAL:HG21	1:A:70:LEU:HD11	1.89	0.55
1:B:12:PRO:HB3	1:B:154:HIS:CD2	2.42	0.55
1:A:212:LEU:HD23	1:A:218:PHE:CE1	2.42	0.55
1:C:226:LEU:O	1:C:230:LEU:HG	2.07	0.55
1:A:162:PRO:HA	1:A:198:GLN:NE2	2.21	0.54
1:B:18:VAL:HG13	1:B:48:THR:HG22	1.89	0.54
1:A:45:GLY:O	1:A:47:THR:HG23	2.07	0.54
1:A:16:ARG:CZ	1:A:251:LEU:HD12	2.37	0.54
1:A:89:THR:HG22	3:A:2019:HOH:O	2.07	0.54
1:A:91:ARG:HB3	1:A:93:GLU:OE1	2.06	0.54
1:C:23:ARG:HG2	1:C:23:ARG:NH2	2.21	0.54
1:B:200:GLN:NE2	1:C:136:ALA:H	2.02	0.53
1:C:16:ARG:HG3	1:C:16:ARG:NH2	2.20	0.53
1:C:32:ARG:HH11	1:C:32:ARG:HG2	1.73	0.53
1:C:270:ARG:O	1:C:273:VAL:HG22	2.08	0.53
1:A:199:LEU:CD1	1:A:206:PRO:HB3	2.37	0.53
1:B:149:SER:CB	1:B:151:LYS:HG2	2.37	0.53
1:B:5:ILE:HG23	1:B:5:ILE:O	2.08	0.53
1:B:149:SER:HB3	1:B:151:LYS:HE2	1.90	0.53
1:B:247:ASN:O	1:B:251:LEU:HB2	2.09	0.53
1:C:195:THR:HG23	3:C:2026:HOH:O	2.07	0.53
1:A:124:LEU:HA	1:A:154:HIS:O	2.09	0.53
1:C:297:VAL:HG22	3:C:2051:HOH:O	2.08	0.53
1:A:70:LEU:O	1:A:70:LEU:HD12	2.09	0.53
1:A:282:ARG:HH12	1:B:102:ASP:HB3	1.72	0.53
1:C:273:VAL:HG23	1:C:274:GLU:OE2	2.09	0.53
1:A:15:SER:O	1:A:17:PHE:N	2.43	0.52
1:C:250:VAL:CA	3:C:2038:HOH:O	2.57	0.52
1:B:5:ILE:HG21	1:B:258:LEU:HD21	1.91	0.52
1:C:213:GLY:O	1:C:216:ARG:HG2	2.08	0.52
1:A:96:ILE:HG22	1:C:283:GLN:HG2	1.92	0.52
1:C:161:THR:HG22	1:C:163:ALA:H	1.75	0.52
1:C:103:ARG:HE	1:C:107:ILE:HD12	1.74	0.52
1:A:212:LEU:HD12	1:A:212:LEU:N	2.24	0.52
1:A:38:ILE:HA	1:A:41:HIS:CD2	2.45	0.52
1:A:57:GLY:O	1:A:58:TYR:HB2	2.11	0.52
1:B:286:PHE:CD1	1:C:98:HIS:HB2	2.45	0.51
1:B:56:GLY:O	1:B:59:GLN:HG3	2.10	0.51
1:C:70:LEU:HD12	1:C:70:LEU:O	2.10	0.51
1:B:32:ARG:HA	1:B:32:ARG:NE	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:ASP:HB2	1:B:83:LYS:O	2.10	0.51
1:A:268:ARG:HH11	2:A:1298:NAP:P2B	2.34	0.51
1:B:283:GLN:O	1:B:287:ARG:HB2	2.11	0.51
1:A:184:SER:OG	1:A:186:VAL:HG22	2.11	0.51
1:C:181:VAL:HB	1:C:190:LEU:HD13	1.93	0.51
1:A:218:PHE:CZ	1:A:243:GLU:HG3	2.46	0.51
1:A:96:ILE:HG12	1:A:97:GLY:N	2.26	0.51
1:B:193:ASP:C	1:B:193:ASP:OD2	2.48	0.51
1:C:192:LEU:HD23	1:C:192:LEU:H	1.75	0.51
1:A:18:VAL:HG23	1:A:260:ILE:HG23	1.91	0.51
1:B:3:GLN:HG2	1:B:4:ARG:O	2.11	0.50
1:C:247:ASN:O	1:C:251:LEU:HB2	2.11	0.50
1:C:271:ALA:O	1:C:274:GLU:HB2	2.11	0.50
1:A:20:GLY:HA3	2:A:1298:NAP:O3D	2.11	0.50
1:B:127:ILE:HG12	1:B:143:PHE:HZ	1.76	0.50
1:A:201:GLN:HA	1:B:171:ARG:NH2	2.25	0.50
1:B:228:ASP:O	1:B:232:VAL:HG23	2.11	0.50
1:B:205:ARG:HD3	3:B:2061:HOH:O	2.10	0.50
1:C:112:GLN:HG3	1:C:116:ASN:ND2	2.23	0.50
1:A:95:VAL:HG22	3:A:2017:HOH:O	2.10	0.50
1:B:79:GLU:HA	1:B:123:ASP:OD2	2.12	0.50
1:A:98:HIS:CE1	1:C:283:GLN:HG3	2.47	0.50
1:B:52:ALA:HB3	1:B:55:TYR:CG	2.47	0.50
1:C:289:ARG:NH2	1:C:295:TYR:O	2.45	0.50
1:A:124:LEU:HG	1:A:126:LEU:HD13	1.94	0.50
1:B:42:LEU:HD13	1:B:77:ARG:CZ	2.42	0.50
1:B:183:ILE:O	1:B:211:CYS:HA	2.12	0.49
1:B:54:ILE:HG22	1:B:58:TYR:CZ	2.47	0.49
1:C:126:LEU:HA	1:C:156:GLY:O	2.11	0.49
1:A:203:ARG:HG2	1:A:203:ARG:HH11	1.77	0.49
1:A:22:TRP:HB2	2:A:1298:NAP:H2D	1.94	0.49
1:B:42:LEU:HD12	1:B:43:ASP:N	2.27	0.49
1:C:107:ILE:O	1:C:111:GLU:HG3	2.12	0.49
1:C:27:TRP:HE3	1:C:29:MET:CE	2.25	0.49
1:C:180:GLN:HE22	2:C:1299:NAP:H2N	1.76	0.49
1:B:161:THR:OG1	1:B:164:GLN:HG3	2.12	0.49
1:A:18:VAL:HG12	1:A:48:THR:HB	1.94	0.49
1:A:54:ILE:HG13	1:A:55:TYR:N	2.27	0.49
1:C:52:ALA:HB2	1:C:83:LYS:HZ3	1.76	0.49
1:B:296:ASP:HB3	3:C:2014:HOH:O	2.11	0.49
1:C:180:GLN:HG2	1:C:209:TRP:HD1	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:VAL:HG12	1:B:273:VAL:HG13	1.95	0.48
1:A:95:VAL:N	1:C:283:GLN:NE2	2.44	0.48
1:B:190:LEU:HD23	3:B:2056:HOH:O	2.13	0.48
1:A:218:PHE:CE2	1:A:243:GLU:HG3	2.48	0.48
1:A:5:ILE:HD11	1:A:205:ARG:HG2	1.95	0.48
1:C:4:ARG:HD3	1:C:13:GLU:OE1	2.14	0.48
1:B:200:GLN:HE22	1:C:136:ALA:N	2.04	0.48
1:A:281:THR:OG1	1:A:284:GLN:HG3	2.13	0.48
1:C:32:ARG:NH1	1:C:32:ARG:HG2	2.28	0.47
1:A:268:ARG:NH1	2:A:1298:NAP:P2B	2.87	0.47
1:C:188:GLN:HE22	1:C:289:ARG:HG3	1.79	0.47
1:C:233:VAL:HG13	1:C:284:GLN:HG2	1.96	0.47
1:C:78:MET:HE2	1:C:80:ILE:HD11	1.97	0.47
1:C:242:ILE:O	1:C:246:VAL:HG23	2.14	0.47
1:A:119:THR:CG2	1:A:120:ASP:N	2.76	0.47
1:A:161:THR:HG21	3:C:2030:HOH:O	2.14	0.47
1:A:212:LEU:HD23	1:A:218:PHE:CZ	2.49	0.47
1:B:42:LEU:HD13	1:B:77:ARG:NH2	2.30	0.47
1:A:32:ARG:HG3	1:A:33:GLN:N	2.30	0.47
1:C:282:ARG:O	1:C:286:PHE:HD1	1.98	0.47
1:A:107:ILE:HD13	1:A:145:HIS:CD2	2.50	0.47
1:A:136:ALA:H	1:C:200:GLN:HE22	1.61	0.46
1:C:162:PRO:HA	1:C:198:GLN:NE2	2.30	0.46
1:C:192:LEU:HD23	1:C:192:LEU:N	2.30	0.46
1:C:86:ILE:HG13	1:C:128:HIS:HB3	1.98	0.46
1:C:91:ARG:HB2	1:C:94:ASN:ND2	2.30	0.46
1:B:199:LEU:HD13	1:B:206:PRO:HB3	1.96	0.46
1:B:216:ARG:HA	1:B:219:ASN:ND2	2.26	0.46
1:B:99:TYR:HB2	1:B:131:ASP:OD1	2.15	0.46
1:B:107:ILE:HD13	1:B:145:HIS:CD2	2.50	0.46
1:C:61:GLU:O	1:C:117:LEU:HD23	2.16	0.46
1:C:237:LEU:O	1:C:238:ASN:HB2	2.15	0.46
1:A:235:GLU:HG3	1:A:236:GLU:N	2.30	0.46
1:C:147:HIS:C	1:C:149:SER:H	2.18	0.46
1:C:7:ILE:HD13	1:C:14:PHE:CD2	2.50	0.46
1:C:250:VAL:CB	3:C:2038:HOH:O	2.45	0.46
1:C:25:MET:CE	1:C:60:CYS:HA	2.46	0.46
1:A:258:LEU:N	1:A:258:LEU:HD23	2.31	0.46
1:B:220:ASP:HB3	1:B:223:PHE:CD1	2.46	0.46
1:A:189:PRO:O	1:A:193:ASP:HB3	2.15	0.46
1:A:80:ILE:HG21	1:A:117:LEU:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:LEU:HD22	1:B:126:LEU:HD12	1.96	0.46
1:B:188:GLN:N	1:B:189:PRO:CD	2.79	0.46
1:C:87:ALA:HB2	1:C:100:ILE:CD1	2.46	0.46
1:B:285:TRP:CE2	1:C:133:LEU:HD21	2.52	0.45
1:C:7:ILE:HG21	1:C:14:PHE:CD2	2.51	0.45
1:A:137:ASP:O	1:A:141:ASP:HB2	2.17	0.45
1:B:147:HIS:HD2	1:B:147:HIS:O	1.99	0.45
1:B:218:PHE:HA	1:B:227:ARG:HD3	1.98	0.45
1:C:51:HIS:O	1:C:82:SER:HA	2.15	0.45
1:A:12:PRO:HD3	1:A:154:HIS:CG	2.52	0.45
1:A:223:PHE:O	1:A:227:ARG:HG3	2.17	0.45
1:A:133:LEU:HA	1:C:254:PRO:HD3	1.99	0.45
1:B:266:ILE:HG13	1:B:266:ILE:O	2.16	0.45
1:B:58:TYR:CE2	1:B:88:THR:OG1	2.70	0.45
1:A:252:ARG:HG3	1:A:252:ARG:HH21	1.80	0.45
1:B:9:PRO:O	1:B:10:GLN:HB2	2.17	0.45
1:A:127:ILE:CG1	1:A:127:ILE:O	2.64	0.45
1:A:199:LEU:HD13	1:A:206:PRO:HB3	1.99	0.45
1:C:121:HIS:CD2	1:C:153:ARG:NE	2.81	0.45
1:C:284:GLN:O	1:C:288:ILE:HG13	2.17	0.45
1:C:2:VAL:HG23	1:C:2:VAL:O	2.16	0.45
1:B:58:TYR:HE2	1:B:88:THR:OG1	2.00	0.44
1:B:105:HIS:HE1	3:B:2029:HOH:O	1.99	0.44
1:B:188:GLN:HB3	1:B:191:LEU:HD22	1.99	0.44
1:C:103:ARG:HH11	1:C:103:ARG:HG3	1.81	0.44
1:A:185:PRO:HG3	1:A:249:TRP:CE3	2.53	0.44
1:B:253:LEU:HD21	1:B:285:TRP:HZ2	1.81	0.44
1:A:179:ASN:HB3	1:A:199:LEU:HD11	2.00	0.44
1:C:84:CYS:SG	1:C:106:ILE:HG23	2.57	0.44
1:A:180:GLN:OE1	2:A:1298:NAP:H2N	2.18	0.44
1:A:205:ARG:N	1:A:205:ARG:HD2	2.32	0.44
1:B:131:ASP:HB3	1:B:134:MET:HB2	2.00	0.44
1:B:190:LEU:N	1:B:190:LEU:HD22	2.33	0.44
1:C:224:GLN:N	1:C:225:PRO:CD	2.80	0.44
1:C:42:LEU:HD13	1:C:77:ARG:CZ	2.48	0.44
1:C:203:ARG:HH11	1:C:203:ARG:HG2	1.83	0.44
1:C:52:ALA:HB2	1:C:83:LYS:NZ	2.31	0.44
1:A:182:GLU:HB2	1:A:209:TRP:CE2	2.52	0.44
1:B:16:ARG:HH22	1:B:256:GLN:HB3	1.83	0.44
1:B:7:ILE:HD11	1:B:177:ALA:HB1	2.00	0.44
1:B:103:ARG:HG2	1:B:103:ARG:NH1	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:ILE:O	1:B:55:TYR:C	2.56	0.43
1:C:7:ILE:HG21	1:C:14:PHE:HD2	1.83	0.43
1:A:228:ASP:O	1:A:232:VAL:HG23	2.17	0.43
1:B:146:LEU:HD13	1:B:152:VAL:HG21	2.00	0.43
1:A:175:THR:HB	3:A:2004:HOH:O	2.18	0.43
1:B:12:PRO:HD3	1:B:154:HIS:CG	2.53	0.43
1:C:182:GLU:HG2	1:C:187:HIS:CD2	2.53	0.43
1:A:217:LEU:HD23	1:A:217:LEU:C	2.38	0.43
1:C:251:LEU:HD13	1:C:259:PRO:HG3	2.00	0.43
1:B:124:LEU:HD22	1:B:126:LEU:CD1	2.49	0.43
1:B:64:PHE:O	1:B:68:LEU:HD13	2.18	0.43
1:B:114:LEU:HA	1:B:114:LEU:HD12	1.87	0.43
1:B:6:THR:HA	1:B:13:GLU:HA	1.99	0.43
1:C:129:ARG:HD3	1:C:130:PRO:HD2	2.01	0.43
1:A:252:ARG:NH2	1:A:252:ARG:HG3	2.34	0.43
1:B:144:LYS:O	1:B:148:GLN:HG3	2.19	0.43
1:B:103:ARG:HH11	1:B:103:ARG:HG2	1.84	0.43
1:B:74:LEU:O	1:B:76:GLU:N	2.52	0.43
1:B:230:LEU:HB3	1:B:242:ILE:HB	2.01	0.43
1:C:140:ALA:HB2	1:C:171:ARG:HB3	2.01	0.43
1:C:83:LYS:HA	1:C:126:LEU:O	2.19	0.43
1:A:271:ALA:O	1:A:274:GLU:HB2	2.19	0.42
1:A:65:GLY:HA3	1:A:116:ASN:O	2.19	0.42
1:A:282:ARG:NH1	1:B:102:ASP:HB3	2.34	0.42
1:B:86:ILE:HG22	1:B:87:ALA:N	2.33	0.42
1:C:129:ARG:CG	1:C:130:PRO:HD2	2.49	0.42
1:C:5:ILE:HG23	1:C:5:ILE:O	2.19	0.42
1:A:161:THR:HG22	1:A:162:PRO:N	2.35	0.42
1:A:266:ILE:O	1:A:270:ARG:HG3	2.18	0.42
1:B:221:ASP:C	1:B:223:PHE:H	2.23	0.42
1:C:129:ARG:HG3	1:C:130:PRO:HD2	2.00	0.42
1:C:35:VAL:HG11	1:C:71:ALA:HB2	2.01	0.42
1:C:25:MET:HG2	1:C:60:CYS:SG	2.60	0.42
1:B:180:GLN:HE21	2:B:1299:NAP:H71N	1.67	0.42
1:C:129:ARG:CD	1:C:130:PRO:HD2	2.50	0.42
1:A:33:GLN:O	1:A:36:SER:N	2.51	0.42
1:C:243:GLU:HG3	1:C:243:GLU:H	1.47	0.42
1:C:30:SER:OG	1:C:33:GLN:HG3	2.19	0.42
1:C:99:TYR:HD2	1:C:131:ASP:OD1	2.02	0.42
1:B:39:GLU:O	1:B:43:ASP:HB2	2.20	0.42
1:C:112:GLN:O	1:C:116:ASN:ND2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:LEU:HB2	2:C:1299:NAP:H52A	2.02	0.42
1:C:88:THR:CG2	1:C:89:THR:N	2.82	0.42
1:B:286:PHE:HZ	1:C:131:ASP:OD1	2.03	0.42
1:A:289:ARG:NH2	1:A:296:ASP:O	2.53	0.42
1:B:44:LEU:HD21	1:B:270:ARG:HB2	2.01	0.42
1:B:28:ASN:ND2	3:B:2010:HOH:O	2.52	0.42
1:C:184:SER:OG	1:C:186:VAL:HG22	2.20	0.42
1:C:51:HIS:O	1:C:83:LYS:NZ	2.47	0.42
1:A:111:GLU:OE2	1:A:151:LYS:HE3	2.19	0.42
1:A:188:GLN:N	1:A:189:PRO:HD2	2.35	0.42
1:A:19:MET:HE2	1:A:261:ILE:HB	2.02	0.42
1:B:202:LEU:O	1:B:203:ARG:HB2	2.20	0.42
1:A:119:THR:CG2	1:A:120:ASP:H	2.30	0.41
1:B:180:GLN:HG3	1:B:207:MET:O	2.19	0.41
1:B:19:MET:HG3	1:B:41:HIS:CD2	2.55	0.41
1:C:8:ALA:HB2	1:C:177:ALA:HA	2.02	0.41
1:C:124:LEU:HD13	1:C:126:LEU:CD1	2.51	0.41
1:C:181:VAL:O	1:C:208:ALA:HA	2.20	0.41
1:B:65:GLY:HA3	1:B:116:ASN:O	2.21	0.41
1:B:5:ILE:HD11	1:B:205:ARG:HG2	2.02	0.41
1:C:233:VAL:HG22	1:C:287:ARG:NH2	2.35	0.41
1:A:283:GLN:HG3	1:B:96:ILE:HG22	2.03	0.41
1:C:6:THR:HG22	1:C:8:ALA:O	2.21	0.41
1:C:88:THR:HG22	1:C:89:THR:N	2.36	0.41
1:C:6:THR:HA	1:C:12:PRO:O	2.20	0.41
1:C:8:ALA:HB1	1:C:9:PRO:CD	2.51	0.41
1:A:75:ARG:NH1	1:A:80:ILE:HD12	2.29	0.41
1:C:52:ALA:HB3	1:C:55:TYR:CD2	2.56	0.41
1:A:267:GLU:H	1:A:267:GLU:CD	2.23	0.41
1:B:32:ARG:HA	1:B:32:ARG:HE	1.83	0.41
1:B:168:LEU:O	1:B:169:GLN:C	2.59	0.41
1:C:40:GLU:HG3	3:C:2006:HOH:O	2.20	0.41
1:C:55:TYR:HB2	3:C:2008:HOH:O	2.20	0.41
1:A:131:ASP:HB3	1:A:134:MET:HB2	2.03	0.41
1:C:218:PHE:CD1	1:C:218:PHE:N	2.88	0.41
1:A:6:THR:HG22	1:A:6:THR:O	2.21	0.40
1:B:4:ARG:HD2	1:B:13:GLU:OE1	2.22	0.40
1:C:203:ARG:HG2	1:C:203:ARG:NH1	2.37	0.40
1:C:259:PRO:HD3	3:C:2039:HOH:O	2.20	0.40
1:B:234:ALA:O	1:B:238:ASN:N	2.52	0.40
1:C:57:GLY:O	1:C:58:TYR:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:GLN:HG3	1:B:116:ASN:HD21	1.87	0.40
1:A:224:GLN:N	1:A:225:PRO:CD	2.84	0.40
1:B:106:ILE:CD1	1:B:139:VAL:HG22	2.51	0.40
1:B:151:LYS:HB3	1:B:151:LYS:NZ	2.35	0.40
1:B:221:ASP:C	1:B:223:PHE:N	2.73	0.40
1:C:27:TRP:HE3	1:C:29:MET:HE1	1.86	0.40
1:C:2:VAL:CG2	1:C:2:VAL:O	2.69	0.40
1:A:22:TRP:CE3	2:A:1298:NAP:H6N	2.56	0.40
1:B:52:ALA:HB3	1:B:55:TYR:CE2	2.57	0.40
1:C:216:ARG:HD2	1:C:220:ASP:OD2	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	296/298 (99%)	266 (90%)	24 (8%)	6 (2%)	9	28
1	B	290/298 (97%)	258 (89%)	26 (9%)	6 (2%)	8	27
1	C	296/298 (99%)	260 (88%)	31 (10%)	5 (2%)	11	34
All	All	882/894 (99%)	784 (89%)	81 (9%)	17 (2%)	9	30

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	ARG
1	A	53	ASP
1	B	75	ARG
1	B	136	ALA
1	C	3	GLN
1	C	28	ASN
1	A	296	ASP

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Mol	Chain	Res	Type
1	B	55	TYR
1	C	53	ASP
1	C	272	ALA
1	C	273	VAL
1	A	128	HIS
1	B	72	PRO
1	A	222	TYR
1	B	5	ILE
1	B	188	GLN
1	A	9	PRO

### 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	252/253 (100%)	236 (94%)	16 (6%)	21	51
1	B	252/253 (100%)	233 (92%)	19 (8%)	16	41
1	C	253/253 (100%)	235 (93%)	18 (7%)	17	44
All	All	757/759 (100%)	704 (93%)	53 (7%)	18	45

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	9	PRO
1	A	19	MET
1	A	42	LEU
1	A	75	ARG
1	A	125	LEU
1	A	126	LEU
1	A	128	HIS
1	A	174	PHE
1	A	180	GLN
1	A	190	LEU
1	A	193	ASP

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Mol	Chain	Res	Type
1	A	205	ARG
1	A	258	LEU
1	A	289	ARG
1	A	293	LEU
1	B	4	ARG
1	B	19	MET
1	B	26	ASP
1	B	42	LEU
1	B	43	ASP
1	B	44	LEU
1	B	75	ARG
1	B	93	GLU
1	B	114	LEU
1	B	124	LEU
1	B	128	HIS
1	B	129	ARG
1	B	151	LYS
1	B	205	ARG
1	B	235	GLU
1	B	238	ASN
1	B	243	GLU
1	B	252	ARG
1	B	287	ARG
1	C	19	MET
1	C	23	ARG
1	C	32	ARG
1	C	41	HIS
1	C	42	LEU
1	C	103	ARG
1	C	104	ASP
1	C	124	LEU
1	C	125	LEU
1	C	174	PHE
1	C	175	THR
1	C	192	LEU
1	C	205	ARG
1	C	218	PHE
1	C	220	ASP
1	C	247	ASN
1	C	284	GLN
1	C	287	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	121	HIS
1	A	128	HIS
1	A	145	HIS
1	A	148	GLN
1	A	154	HIS
1	A	169	GLN
1	A	188	GLN
1	A	198	GLN
1	A	200	GLN
1	A	201	GLN
1	A	219	ASN
1	B	28	ASN
1	B	33	GLN
1	B	105	HIS
1	B	112	GLN
1	B	116	ASN
1	B	145	HIS
1	B	147	HIS
1	B	148	GLN
1	B	154	HIS
1	B	169	GLN
1	B	180	GLN
1	B	198	GLN
1	B	200	GLN
1	B	201	GLN
1	B	219	ASN
1	B	244	GLN
1	C	3	GLN
1	C	10	GLN
1	C	28	ASN
1	C	59	GLN
1	C	105	HIS
1	C	116	ASN
1	C	121	HIS
1	C	145	HIS
1	C	148	GLN
1	C	154	HIS
1	C	169	GLN
1	C	180	GLN
1	C	187	HIS
1	C	188	GLN

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Mol	Chain	Res	Type
1	C	198	GLN
1	C	200	GLN
1	C	201	GLN
1	C	238	ASN
1	C	244	GLN
1	C	247	ASN
1	C	283	GLN

### 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 ⓘ

3 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	NAP	A	1298	-	44,52,52	1.83	10 (22%)	51,80,80	2.71	20 (39%)
2	NAP	B	1299	-	44,52,52	1.84	11 (25%)	51,80,80	2.69	21 (41%)
2	NAP	C	1299	-	44,52,52	1.79	11 (25%)	51,80,80	2.71	19 (37%)

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	NAP	A	1298	-	-	0/27/67/67	0/5/5/5
2	NAP	B	1299	-	-	0/27/67/67	0/5/5/5
2	NAP	C	1299	-	-	0/27/67/67	0/5/5/5

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1299	NAP	C2D-C1D	-3.83	1.47	1.53
2	A	1298	NAP	PA-O1A	-2.60	1.41	1.50
2	C	1299	NAP	C2D-C1D	-2.51	1.49	1.53
2	C	1299	NAP	PA-O1A	-2.50	1.41	1.50
2	A	1298	NAP	C2D-C1D	-2.49	1.49	1.53
2	B	1299	NAP	PA-O1A	-2.28	1.42	1.50
2	C	1299	NAP	C6N-C5N	2.04	1.43	1.38
2	B	1299	NAP	C3N-C7N	2.05	1.53	1.50
2	B	1299	NAP	C6N-C5N	2.05	1.43	1.38
2	B	1299	NAP	C5N-C4N	2.06	1.42	1.38
2	C	1299	NAP	C5N-C4N	2.14	1.43	1.38
2	C	1299	NAP	C3N-C7N	2.19	1.53	1.50
2	A	1298	NAP	P2B-O2B	2.30	1.63	1.59
2	B	1299	NAP	C4A-N3A	2.31	1.39	1.35
2	C	1299	NAP	C4A-N3A	2.45	1.39	1.35
2	A	1298	NAP	C4A-N3A	2.68	1.39	1.35
2	A	1298	NAP	C3N-C7N	2.75	1.54	1.50
2	B	1299	NAP	C6N-N1N	3.13	1.43	1.35
2	A	1298	NAP	C6N-N1N	3.26	1.43	1.35
2	C	1299	NAP	C6N-N1N	3.47	1.44	1.35
2	B	1299	NAP	O4D-C1D	3.70	1.46	1.41
2	C	1299	NAP	C2A-N1A	3.82	1.41	1.33
2	C	1299	NAP	O4D-C1D	3.88	1.46	1.41
2	A	1298	NAP	C2A-N1A	3.89	1.41	1.33
2	A	1298	NAP	O4D-C1D	3.89	1.46	1.41
2	B	1299	NAP	C2A-N1A	3.91	1.41	1.33
2	A	1298	NAP	C2A-N3A	4.10	1.39	1.32
2	B	1299	NAP	C2A-N3A	4.17	1.39	1.32
2	C	1299	NAP	C2A-N3A	4.27	1.39	1.32
2	C	1299	NAP	C4N-C3N	4.82	1.47	1.39
2	A	1298	NAP	C4N-C3N	4.89	1.47	1.39
2	B	1299	NAP	C4N-C3N	5.18	1.47	1.39

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1299	NAP	O3X-P2B-O1X	-7.73	80.25	110.50
2	B	1299	NAP	O3X-P2B-O1X	-7.68	80.45	110.50
2	A	1298	NAP	O5D-PN-O1N	-7.53	78.86	109.25
2	C	1299	NAP	O5D-PN-O1N	-7.48	79.07	109.25
2	B	1299	NAP	O5D-PN-O1N	-7.33	79.67	109.25
2	A	1298	NAP	O3X-P2B-O1X	-7.22	82.24	110.50
2	C	1299	NAP	N3A-C2A-N1A	-6.75	122.98	128.86
2	A	1298	NAP	N3A-C2A-N1A	-6.71	123.01	128.86
2	B	1299	NAP	N3A-C2A-N1A	-6.63	123.08	128.86
2	B	1299	NAP	O3X-P2B-O2B	-5.90	79.19	106.00
2	C	1299	NAP	O3X-P2B-O2B	-5.88	79.28	106.00
2	A	1298	NAP	O3X-P2B-O2B	-5.80	79.64	106.00
2	C	1299	NAP	O2N-PN-O5D	-4.07	88.94	108.14
2	A	1298	NAP	O2N-PN-O5D	-4.05	89.01	108.14
2	A	1298	NAP	O3X-P2B-O2X	-4.01	91.43	107.61
2	B	1299	NAP	O2N-PN-O5D	-3.99	89.28	108.14
2	B	1299	NAP	O3X-P2B-O2X	-3.76	92.45	107.61
2	C	1299	NAP	O3X-P2B-O2X	-3.67	92.80	107.61
2	A	1298	NAP	O7N-C7N-C3N	-2.82	116.33	119.62
2	C	1299	NAP	O7N-C7N-C3N	-2.81	116.34	119.62
2	B	1299	NAP	O7N-C7N-C3N	-2.71	116.46	119.62
2	A	1298	NAP	O4B-C4B-C5B	-2.55	100.81	109.40
2	C	1299	NAP	C1B-N9A-C4A	-2.43	122.44	126.64
2	A	1298	NAP	O3B-C3B-C4B	-2.42	104.03	111.09
2	B	1299	NAP	O4B-C4B-C5B	-2.36	101.42	109.40
2	B	1299	NAP	C4D-O4D-C1D	-2.29	107.33	109.77
2	A	1298	NAP	C1B-N9A-C4A	-2.19	122.84	126.64
2	C	1299	NAP	O4B-C4B-C5B	-2.10	102.32	109.40
2	B	1299	NAP	C5N-C4N-C3N	-2.09	117.89	120.35
2	B	1299	NAP	O3B-C3B-C4B	-2.07	105.03	111.09
2	A	1298	NAP	C3N-C2N-N1N	-2.06	118.35	120.43
2	C	1299	NAP	C5N-C4N-C3N	-2.04	117.94	120.35
2	B	1299	NAP	C1B-N9A-C4A	-2.04	123.11	126.64
2	B	1299	NAP	C2B-C3B-C4B	2.01	106.53	101.95
2	B	1299	NAP	O2N-PN-O1N	2.02	122.71	112.28
2	C	1299	NAP	O2N-PN-O1N	2.14	123.35	112.28
2	A	1298	NAP	O2N-PN-O1N	2.15	123.41	112.28
2	A	1298	NAP	C2B-C3B-C4B	2.18	106.91	101.95
2	C	1299	NAP	C2B-C3B-C4B	2.37	107.33	101.95
2	B	1299	NAP	O2X-P2B-O1X	2.49	120.24	110.50
2	C	1299	NAP	O2X-P2B-O1X	2.51	120.33	110.50
2	A	1298	NAP	O2A-PA-O1A	2.54	125.42	112.28
2	A	1298	NAP	C2N-C3N-C4N	2.56	121.18	118.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1298	NAP	O2X-P2B-O1X	2.56	120.51	110.50
2	C	1299	NAP	O2A-PA-O1A	2.63	125.91	112.28
2	B	1299	NAP	C2N-C3N-C4N	2.74	121.39	118.26
2	B	1299	NAP	O2A-PA-O1A	2.75	126.51	112.28
2	C	1299	NAP	C2N-C3N-C4N	2.82	121.47	118.26
2	B	1299	NAP	C3N-C7N-N7N	3.12	121.34	117.77
2	A	1298	NAP	C5A-C6A-N6A	3.17	126.93	120.47
2	C	1299	NAP	C5A-C6A-N6A	3.26	127.12	120.47
2	C	1299	NAP	C3N-C7N-N7N	3.29	121.53	117.77
2	B	1299	NAP	C5A-C6A-N6A	3.32	127.24	120.47
2	A	1298	NAP	C3N-C7N-N7N	3.69	121.99	117.77
2	A	1298	NAP	O2B-P2B-O1X	4.36	126.36	109.26
2	C	1299	NAP	C4B-O4B-C1B	4.67	114.74	109.77
2	B	1299	NAP	C4B-O4B-C1B	4.70	114.77	109.77
2	B	1299	NAP	O2B-P2B-O1X	4.74	127.85	109.26
2	C	1299	NAP	O2B-P2B-O1X	4.80	128.09	109.26
2	A	1298	NAP	C4B-O4B-C1B	4.87	114.95	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1298	NAP	8	0
2	B	1299	NAP	7	0
2	C	1299	NAP	5	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.