



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:05 AM GMT

PDB ID : 2PD3  
Title : Crystal Structure of the Helicobacter pylori Enoyl-Acyl Carrier Protein Reductase in Complex with Hydroxydiphenyl Ether Compounds, Triclosan and Diclosan  
Authors : Lee, H.H.; Moon, J.H.; Suh, S.W.  
Deposited on : 2007-03-31  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

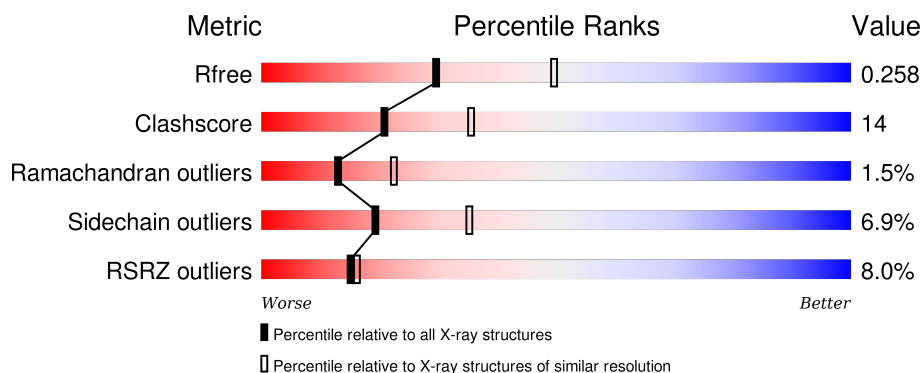
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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.

Mol	Chain	Length	Quality of chain
1	A	275	<div> <div>9%</div> <div>69% 25% 5%</div> </div>
1	B	275	<div> <div>9%</div> <div>69% 28% .</div> </div>
1	C	275	<div> <div>7%</div> <div>69% 27% .</div> </div>
1	D	275	<div> <div>8%</div> <div>67% 29% .</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8866 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 Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2102	1341	354	399	8			
1	B	274	Total	C	N	O	S	0	0	0
			2102	1341	354	399	8			
1	C	274	Total	C	N	O	S	0	0	0
			2102	1341	354	399	8			
1	D	274	Total	C	N	O	S	0	0	0
			2102	1341	354	399	8			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



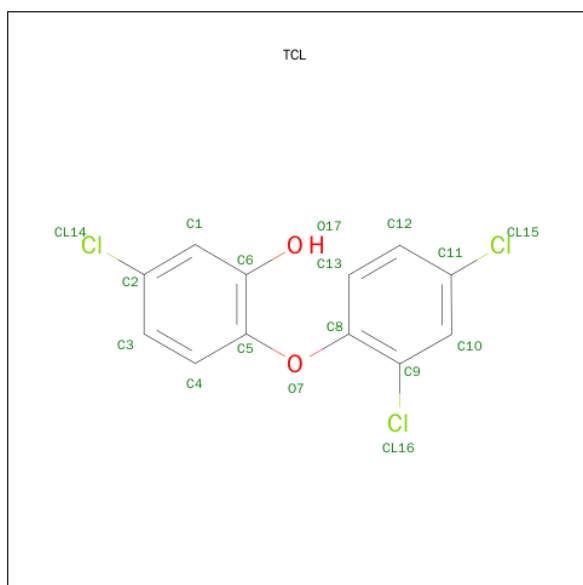
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is TRICLOSAN (three-letter code: TCL) (formula:  $C_{12}H_7Cl_3O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	O	0	0
			17	12	3	2		
3	B	1	Total	C	Cl	O	0	0
			17	12	3	2		
3	C	1	Total	C	Cl	O	0	0
			17	12	3	2		
3	D	1	Total	C	Cl	O	0	0
			17	12	3	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total	O	0	0
			40	40		
4	B	56	Total	O	0	0
			56	56		
4	C	59	Total	O	0	0
			59	59		

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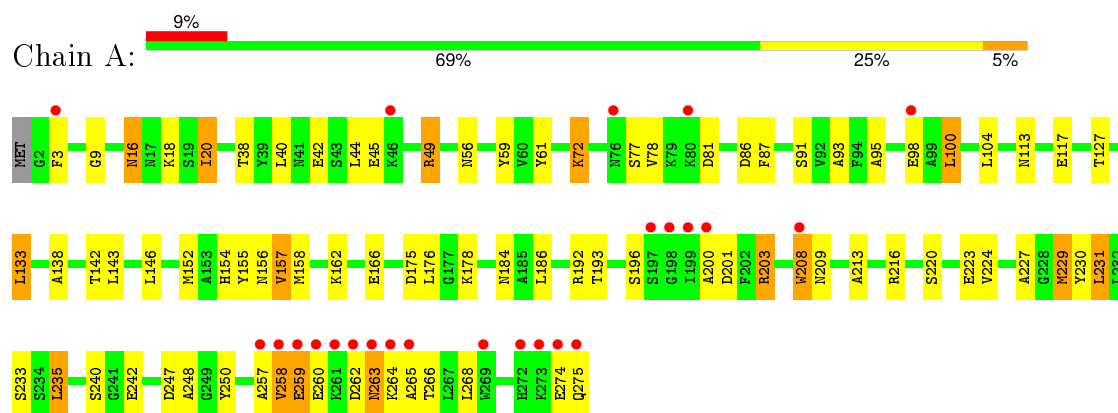
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	59	Total	O	0	0
			59	59		

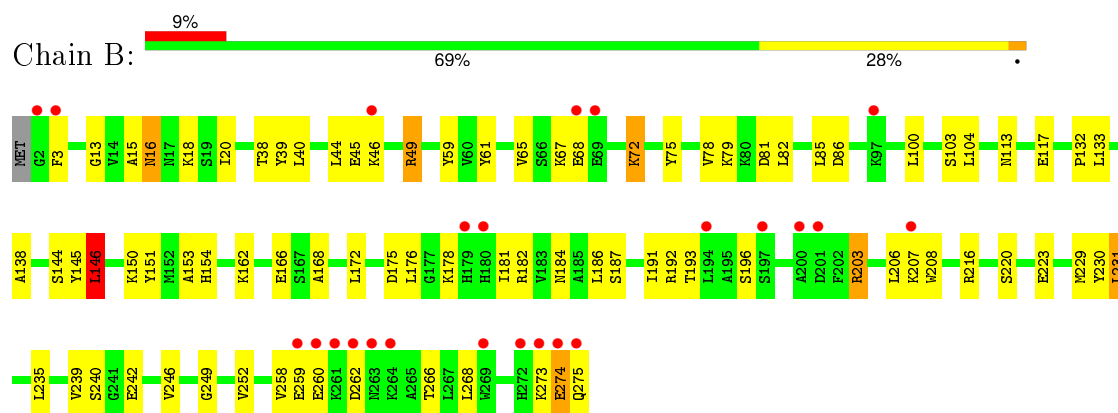
### 3 Residue-property plots

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.

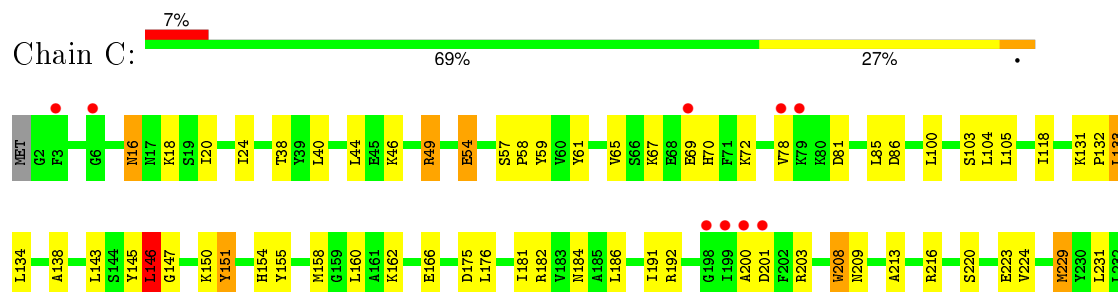
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

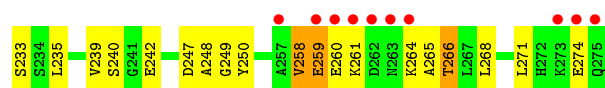


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

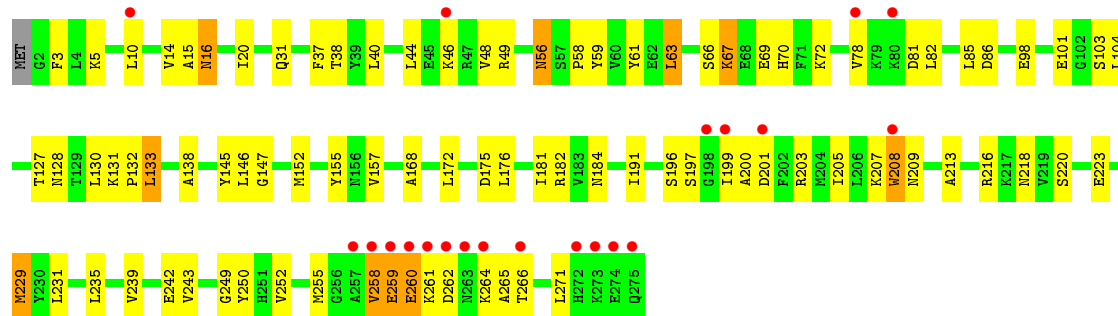


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]





- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.35Å 94.91Å 75.38Å 90.00° 106.21° 90.00°	Depositor
Resolution (Å)	19.82 – 2.50 19.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (19.82-2.50) 99.1 (19.82-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.37 (at 2.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.233 , 0.259 0.231 , 0.258	Depositor DCC
$R_{free}$ test set	3400 reflections (9.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.1	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.5	EDS
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	5 of 34060 reflections (0.015%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8866	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	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 74.38 % 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 1.5457e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.40	0/2140	0.64	1/2890 (0.0%)
1	B	0.41	0/2140	0.65	1/2890 (0.0%)
1	C	0.41	0/2140	0.65	0/2890
1	D	0.41	0/2140	0.64	0/2890
All	All	0.41	0/8560	0.64	2/11560 (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	259	GLU	N-CA-C	5.19	125.02	111.00
1	B	206	LEU	CA-CB-CG	5.08	126.99	115.30

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	2102	0	2125	65	0
1	B	2102	0	2125	61	0
1	C	2102	0	2125	65	0
1	D	2102	0	2125	66	0
2	A	44	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	44	0	26	6	0
2	C	44	0	26	2	0
2	D	44	0	26	4	0
3	A	17	0	6	2	0
3	B	17	0	6	2	0
3	C	17	0	6	3	0
3	D	17	0	6	2	0
4	A	40	0	0	3	0
4	B	56	0	0	3	0
4	C	59	0	0	2	0
4	D	59	0	0	10	0
All	All	8866	0	8628	249	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 14.

All (249) 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:B:216:ARG:HH22	1:B:260:GLU:HB2	1.18	1.05
1:D:184:ASN:HD22	1:D:242:GLU:H	1.12	0.97
1:C:216:ARG:HH22	1:C:260:GLU:HB2	1.28	0.96
1:B:184:ASN:HD22	1:B:242:GLU:H	1.15	0.92
1:C:100:LEU:HD21	3:C:4414:TCL:H121	1.56	0.87
1:B:184:ASN:ND2	1:B:242:GLU:H	1.71	0.87
1:D:78:VAL:CG2	1:D:133:LEU:HD23	2.08	0.84
1:A:40:LEU:HB3	1:A:44:LEU:HD12	1.60	0.84
1:D:184:ASN:ND2	1:D:242:GLU:H	1.77	0.82
1:A:192:ARG:HD3	1:A:203:ARG:HH12	1.43	0.81
1:A:184:ASN:HD21	1:A:240:SER:HA	1.45	0.79
1:A:157:VAL:HB	4:A:2419:HOH:O	1.82	0.78
1:D:203:ARG:O	1:D:207:LYS:HG3	1.84	0.77
1:A:184:ASN:HD22	1:A:242:GLU:H	1.34	0.76
1:A:192:ARG:HG3	1:A:192:ARG:HH11	1.52	0.73
1:D:216:ARG:HH22	1:D:260:GLU:HB2	1.50	0.73
1:C:184:ASN:HD22	1:C:242:GLU:H	1.37	0.72
1:C:16:ASN:ND2	1:C:18:LYS:H	1.88	0.72
1:C:16:ASN:C	1:C:16:ASN:HD22	1.93	0.70
1:C:40:LEU:HB3	1:C:44:LEU:HD12	1.74	0.70
1:B:100:LEU:HD21	3:B:3414:TCL:H121	1.73	0.69
1:A:258:VAL:HG23	1:A:266:THR:O	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ARG:O	1:B:207:LYS:HG3	1.93	0.68
1:A:16:ASN:ND2	1:A:18:LYS:H	1.91	0.68
1:B:113:ASN:O	1:B:117:GLU:HG3	1.94	0.68
1:A:208:TRP:HZ3	1:A:250:TYR:HH	1.39	0.68
1:B:216:ARG:NH2	1:B:260:GLU:HB2	2.01	0.68
1:D:157:VAL:HB	4:D:5447:HOH:O	1.94	0.67
1:B:59:TYR:CE2	1:B:82:LEU:HD21	2.30	0.66
2:A:1780:NAD:O2N	2:A:1780:NAD:H2N	1.95	0.66
1:D:78:VAL:HG23	1:D:133:LEU:HD23	1.77	0.66
1:A:229:MET:CE	1:A:233:SER:HB3	2.26	0.66
1:C:184:ASN:ND2	1:C:242:GLU:H	1.93	0.66
1:D:14:VAL:HG23	4:D:5418:HOH:O	1.95	0.65
1:A:220:SER:OG	1:A:223:GLU:HG3	1.96	0.65
1:D:127:THR:HG21	1:D:176:LEU:HD11	1.76	0.65
1:A:16:ASN:C	1:A:16:ASN:HD22	1.97	0.65
1:B:184:ASN:HD22	1:B:242:GLU:N	1.91	0.65
1:C:86:ASP:O	1:C:138:ALA:HA	1.97	0.64
1:A:216:ARG:HH22	1:A:260:GLU:HB2	1.63	0.64
1:B:40:LEU:HB3	1:B:44:LEU:HD12	1.80	0.64
1:D:184:ASN:HD22	1:D:242:GLU:N	1.92	0.63
1:D:131:LYS:HB3	1:D:132:PRO:HD3	1.80	0.63
1:D:258:VAL:HG23	1:D:266:THR:O	1.98	0.63
1:D:38:THR:HA	1:D:61:TYR:O	1.99	0.61
1:B:178:LYS:HD3	1:D:259:GLU:CG	2.30	0.61
1:A:192:ARG:HG3	1:A:192:ARG:NH1	2.15	0.61
1:B:162:LYS:O	1:B:166:GLU:HG3	2.00	0.61
1:D:208:TRP:HZ3	1:D:250:TYR:HH	1.47	0.60
1:D:249:GLY:O	1:D:252:VAL:HG22	2.01	0.60
1:A:184:ASN:HD21	1:A:240:SER:CA	2.15	0.60
1:A:162:LYS:O	1:A:166:GLU:HG3	2.02	0.60
1:A:208:TRP:HZ3	1:A:250:TYR:OH	1.84	0.60
1:D:203:ARG:NH2	4:D:5443:HOH:O	2.34	0.59
1:D:66:SER:OG	1:D:67:LYS:HD2	2.02	0.59
1:B:16:ASN:C	1:B:16:ASN:HD22	2.06	0.59
1:B:146:LEU:HD22	1:B:150:LYS:HB2	1.85	0.59
1:B:45:GLU:OE1	1:B:49:ARG:NH1	2.35	0.58
1:A:72:LYS:HA	1:A:72:LYS:HE3	1.85	0.58
1:C:216:ARG:NH2	1:C:260:GLU:HB2	2.09	0.57
1:A:86:ASP:O	1:A:138:ALA:HA	2.03	0.57
1:B:78:VAL:HG23	1:B:133:LEU:HD23	1.84	0.57
1:C:162:LYS:O	1:C:166:GLU:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:MET:CE	4:D:5416:HOH:O	2.51	0.57
1:C:20:ILE:HG23	1:C:224:VAL:HG11	1.86	0.57
1:B:59:TYR:CE1	1:B:81:ASP:HB3	2.39	0.57
1:D:266:THR:HG23	4:D:5432:HOH:O	2.04	0.57
1:C:16:ASN:HD22	1:C:18:LYS:H	1.52	0.56
1:D:20:ILE:HD11	1:D:191:ILE:HG21	1.87	0.56
1:A:98:GLU:CD	1:A:98:GLU:H	2.09	0.56
1:C:18:LYS:HD3	4:C:4443:HOH:O	2.04	0.56
1:A:152:MET:CE	4:A:2421:HOH:O	2.53	0.56
1:C:38:THR:HA	1:C:61:TYR:O	2.06	0.56
1:B:75:TYR:HE2	1:B:132:PRO:HG2	1.70	0.56
1:A:230:TYR:CD1	1:A:231:LEU:HD13	2.40	0.56
1:A:127:THR:HG21	1:A:176:LEU:HD11	1.88	0.56
1:A:229:MET:HE3	1:A:233:SER:HB3	1.88	0.55
1:B:38:THR:HA	1:B:61:TYR:O	2.05	0.55
1:A:152:MET:HE2	4:A:2421:HOH:O	2.05	0.55
1:C:247:ASP:O	1:C:248:ALA:HB3	2.07	0.55
1:B:273:LYS:O	1:B:274:GLU:HB2	2.05	0.55
1:D:208:TRP:HZ3	1:D:250:TYR:OH	1.90	0.55
1:D:44:LEU:O	1:D:48:VAL:HG23	2.07	0.55
1:A:216:ARG:HH22	1:A:260:GLU:CB	2.20	0.54
1:D:98:GLU:H	1:D:98:GLU:CD	2.11	0.54
1:A:184:ASN:ND2	1:A:240:SER:HA	2.17	0.54
1:C:18:LYS:HE3	4:C:4451:HOH:O	2.07	0.54
1:B:178:LYS:HD3	1:D:259:GLU:HG3	1.89	0.54
1:B:220:SER:OG	1:B:223:GLU:HG3	2.08	0.53
1:B:262:ASP:OD1	1:C:154:HIS:CE1	2.61	0.53
1:B:178:LYS:HD3	1:D:259:GLU:HG2	1.91	0.53
1:A:155:TYR:CZ	1:A:158:MET:HG3	2.43	0.53
1:D:220:SER:OG	1:D:223:GLU:HG3	2.09	0.53
1:A:38:THR:HA	1:A:61:TYR:O	2.08	0.53
1:C:131:LYS:HB3	1:C:132:PRO:HD3	1.90	0.52
1:A:61:TYR:CE2	1:A:78:VAL:HG12	2.44	0.52
1:A:154:HIS:CE1	1:D:262:ASP:OD1	2.63	0.52
1:B:86:ASP:O	1:B:138:ALA:HA	2.10	0.51
1:B:20:ILE:HD11	1:B:191:ILE:HG21	1.91	0.51
1:C:143:LEU:O	1:C:162:LYS:NZ	2.43	0.51
1:C:220:SER:OG	1:C:223:GLU:HG3	2.11	0.51
1:D:176:LEU:HB3	1:D:181:ILE:HB	1.91	0.51
1:A:258:VAL:HG23	1:A:266:THR:C	2.30	0.51
1:A:230:TYR:HD1	1:A:231:LEU:HD13	1.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:69:GLU:HG2	1:C:70:HIS:N	2.26	0.51
1:C:20:ILE:O	1:C:24:ILE:HG13	2.10	0.51
1:C:155:TYR:CZ	1:C:158:MET:HG3	2.46	0.51
1:B:176:LEU:HB3	1:B:181:ILE:HB	1.92	0.51
1:C:100:LEU:CD2	3:C:4414:TCL:H121	2.37	0.50
1:C:175:ASP:OD2	1:D:103:SER:HB2	2.12	0.50
1:C:49:ARG:HG3	1:C:49:ARG:HH11	1.77	0.50
1:D:31:GLN:HG3	1:D:229:MET:SD	2.51	0.50
1:A:61:TYR:CE1	1:A:77:SER:HB3	2.47	0.50
1:C:131:LYS:HA	1:C:134:LEU:HD12	1.95	0.49
1:B:100:LEU:CD2	3:B:3414:TCL:H121	2.42	0.49
1:C:146:LEU:HD22	1:C:150:LYS:HB2	1.95	0.49
1:C:20:ILE:HD11	1:C:191:ILE:HG21	1.95	0.48
1:C:151:TYR:CD1	1:C:151:TYR:C	2.86	0.48
1:D:218:ASN:HB3	4:D:5422:HOH:O	2.13	0.48
1:A:20:ILE:HG23	1:A:224:VAL:HG11	1.96	0.48
1:A:186:LEU:HD11	1:A:227:ALA:HB3	1.95	0.48
1:A:3:PHE:CE2	1:A:235:LEU:HD11	2.48	0.48
1:D:59:TYR:CD1	1:D:81:ASP:HB3	2.49	0.48
1:B:3:PHE:CG	1:B:235:LEU:HD13	2.49	0.47
1:B:46:LYS:HD3	1:B:46:LYS:O	2.14	0.47
1:C:258:VAL:HG23	1:C:266:THR:O	2.14	0.47
1:B:216:ARG:NH2	1:B:268:LEU:HD21	2.30	0.47
1:A:9:GLY:HA2	1:A:87:PHE:O	2.13	0.47
2:D:4780:NAD:H52N	3:D:5414:TCL:CL16	2.52	0.47
1:B:260:GLU:HA	1:B:260:GLU:OE1	2.14	0.47
1:B:72:LYS:HA	1:B:72:LYS:HE3	1.96	0.47
1:D:63:LEU:HD23	1:D:70:HIS:HB3	1.97	0.47
1:D:127:THR:CG2	1:D:176:LEU:HD11	2.44	0.47
1:B:186:LEU:HD21	1:B:246:VAL:CG2	2.44	0.47
1:D:152:MET:HE1	4:D:5416:HOH:O	2.13	0.47
1:C:160:LEU:HD11	1:D:168:ALA:HA	1.96	0.47
1:D:3:PHE:CD2	1:D:235:LEU:HD13	2.50	0.46
1:B:68:GLU:CD	4:B:3440:HOH:O	2.53	0.46
1:D:16:ASN:ND2	4:D:5427:HOH:O	2.23	0.46
1:A:113:ASN:O	1:A:117:GLU:HG3	2.15	0.46
1:D:264:LYS:O	1:D:265:ALA:HB3	2.15	0.46
1:A:193:THR:H	1:A:196:SER:HG	1.60	0.46
1:D:197:SER:HA	4:D:5443:HOH:O	2.15	0.46
1:A:100:LEU:HD13	1:A:154:HIS:O	2.15	0.46
1:C:49:ARG:HH11	1:C:49:ARG:CG	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:ARG:HH11	1:B:192:ARG:HG3	1.81	0.46
1:B:182:ARG:HD2	1:B:239:VAL:O	2.16	0.46
1:D:200:ALA:O	1:D:201:ASP:CG	2.53	0.46
1:C:16:ASN:C	1:C:16:ASN:ND2	2.66	0.46
1:B:266:THR:HG23	4:B:3465:HOH:O	2.16	0.46
1:A:142:THR:C	1:A:143:LEU:HD22	2.35	0.46
2:D:4780:NAD:H2N	2:D:4780:NAD:O2N	2.15	0.45
1:B:249:GLY:O	1:B:252:VAL:HG22	2.15	0.45
1:B:15:ALA:HB2	2:B:2780:NAD:O3B	2.16	0.45
1:A:209:ASN:O	1:A:213:ALA:HB2	2.16	0.45
1:A:262:ASP:O	1:A:263:ASN:C	2.54	0.45
2:C:3780:NAD:H2D	3:C:4414:TCL:O7	2.16	0.45
1:B:59:TYR:CD1	1:B:81:ASP:HB3	2.51	0.45
1:C:261:LYS:N	1:C:261:LYS:HD2	2.32	0.45
1:A:78:VAL:CG2	1:A:133:LEU:HD23	2.47	0.45
1:C:208:TRP:HZ3	1:C:250:TYR:OH	2.00	0.45
1:D:5:LYS:HD2	4:D:5464:HOH:O	2.17	0.45
1:B:144:SER:OG	1:B:145:TYR:N	2.50	0.45
1:A:155:TYR:O	1:A:156:ASN:C	2.54	0.44
1:B:175:ASP:O	1:B:178:LYS:HE2	2.18	0.44
1:C:176:LEU:HB3	1:C:181:ILE:HB	2.00	0.44
2:B:2780:NAD:H51N	2:B:2780:NAD:H52A	1.99	0.44
1:B:39:TYR:CD2	1:B:45:GLU:HB2	2.52	0.44
1:A:231:LEU:HA	1:A:231:LEU:HD12	1.85	0.44
1:D:69:GLU:HG2	1:D:70:HIS:N	2.33	0.44
1:C:264:LYS:O	1:C:264:LYS:HG3	2.18	0.44
1:C:54:GLU:O	1:C:54:GLU:HG2	2.17	0.44
1:B:230:TYR:CD1	1:B:231:LEU:HD13	2.53	0.44
1:A:208:TRP:CZ3	1:D:255:MET:HA	2.53	0.44
1:B:153:ALA:HB1	1:C:258:VAL:HG11	2.00	0.44
1:D:147:GLY:HA3	1:D:155:TYR:CD2	2.53	0.44
1:B:151:TYR:C	1:B:151:TYR:CD1	2.91	0.44
1:C:65:VAL:HG21	1:C:118:ILE:HG23	2.00	0.44
1:D:168:ALA:O	1:D:172:LEU:HG	2.17	0.43
1:D:40:LEU:HB3	1:D:44:LEU:HD12	2.00	0.43
1:C:78:VAL:HG23	1:C:133:LEU:HD23	2.00	0.43
1:D:145:TYR:OH	1:D:152:MET:CE	2.66	0.43
1:A:175:ASP:CG	1:B:103:SER:HB2	2.39	0.43
1:B:100:LEU:O	1:B:154:HIS:HB3	2.19	0.43
1:A:264:LYS:O	1:A:265:ALA:HB3	2.18	0.43
1:B:65:VAL:HG22	2:B:2780:NAD:N1A	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:ASP:O	1:A:248:ALA:HB3	2.18	0.43
1:B:16:ASN:ND2	4:B:3427:HOH:O	2.34	0.43
1:D:98:GLU:O	1:D:101:GLU:HG2	2.18	0.43
1:B:186:LEU:C	1:B:186:LEU:HD23	2.39	0.43
1:D:15:ALA:HB2	2:D:4780:NAD:O3B	2.19	0.43
1:C:105:LEU:HB3	1:D:128:ASN:HD22	1.83	0.43
1:A:95:ALA:N	3:A:2414:TCL:CL15	2.88	0.43
1:A:178:LYS:HD3	1:C:259:GLU:CG	2.48	0.43
1:A:45:GLU:OE1	1:A:49:ARG:NH1	2.52	0.43
1:B:193:THR:H	1:B:196:SER:HG	1.65	0.43
1:A:184:ASN:ND2	1:A:242:GLU:H	2.07	0.42
1:D:59:TYR:CE2	1:D:82:LEU:HD21	2.54	0.42
1:C:182:ARG:HD2	1:C:239:VAL:O	2.19	0.42
1:D:205:ILE:HG12	1:D:205:ILE:H	1.73	0.42
1:D:10:LEU:CD2	1:D:130:LEU:HD11	2.50	0.42
1:C:261:LYS:HG2	1:C:271:LEU:CD1	2.49	0.42
1:D:14:VAL:HG21	1:D:37:PHE:HD1	1.85	0.42
1:B:20:ILE:HG12	2:B:2780:NAD:O1N	2.19	0.42
1:D:261:LYS:HG2	1:D:271:LEU:CD1	2.50	0.42
1:D:196:SER:HA	1:D:199:ILE:HD12	2.02	0.42
1:C:200:ALA:O	1:C:201:ASP:CG	2.58	0.42
1:C:145:TYR:O	1:C:147:GLY:N	2.53	0.42
1:D:182:ARG:HD2	1:D:239:VAL:O	2.19	0.42
2:D:4780:NAD:H2D	3:D:5414:TCL:O7	2.19	0.42
1:C:261:LYS:O	1:C:265:ALA:HA	2.19	0.42
1:A:257:ALA:CB	1:A:268:LEU:HD22	2.51	0.41
1:D:264:LYS:HD2	1:D:264:LYS:HA	1.75	0.41
1:C:105:LEU:HB3	1:D:128:ASN:ND2	2.35	0.41
1:C:65:VAL:CG2	1:C:118:ILE:HG23	2.50	0.41
1:A:16:ASN:C	1:A:16:ASN:ND2	2.68	0.41
1:D:86:ASP:O	1:D:138:ALA:HA	2.19	0.41
1:B:231:LEU:HA	1:B:231:LEU:HD12	1.85	0.41
1:C:59:TYR:CD1	1:C:81:ASP:HB3	2.56	0.41
1:B:13:GLY:O	2:B:2780:NAD:H4B	2.20	0.41
1:A:59:TYR:CD1	1:A:81:ASP:HB3	2.56	0.41
1:C:192:ARG:HG3	1:C:192:ARG:HH11	1.86	0.41
1:C:209:ASN:O	1:C:213:ALA:HB2	2.20	0.41
1:C:235:LEU:HA	1:C:235:LEU:HD12	1.94	0.41
1:B:79:LYS:HB2	1:B:133:LEU:HD21	2.03	0.41
1:C:233:SER:OG	1:C:235:LEU:HB2	2.20	0.41
1:A:200:ALA:O	1:A:201:ASP:CG	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:216:ARG:NH2	1:C:268:LEU:HD21	2.36	0.41
1:D:209:ASN:O	1:D:213:ALA:HB2	2.20	0.41
1:C:65:VAL:HG22	2:C:3780:NAD:N1A	2.35	0.41
1:A:240:SER:O	1:C:249:GLY:HA2	2.21	0.41
1:A:216:ARG:HH21	1:A:268:LEU:HD21	1.85	0.41
1:B:16:ASN:ND2	1:B:16:ASN:C	2.73	0.41
1:B:16:ASN:ND2	1:B:18:LYS:H	2.18	0.41
1:B:186:LEU:HD23	1:B:187:SER:N	2.35	0.41
1:A:93:ALA:O	3:A:2414:TCL:H101	2.21	0.41
1:D:56:ASN:HD22	1:D:56:ASN:HA	1.66	0.41
1:C:57:SER:HA	1:C:58:PRO:HD3	1.89	0.41
1:A:233:SER:OG	1:A:235:LEU:HB2	2.21	0.41
1:A:9:GLY:HA3	1:A:87:PHE:CE1	2.55	0.40
2:B:2780:NAD:H2N	2:B:2780:NAD:O2N	2.22	0.40
1:C:229:MET:HE2	1:C:233:SER:HB3	2.03	0.40
1:C:103:SER:HB2	1:D:175:ASP:CG	2.41	0.40
1:C:260:GLU:HA	1:C:260:GLU:OE1	2.21	0.40
1:C:49:ARG:NH1	1:C:49:ARG:CG	2.84	0.40
1:C:186:LEU:C	1:C:186:LEU:HD23	2.41	0.40
1:D:242:GLU:HG3	1:D:243:VAL:N	2.36	0.40
1:A:127:THR:CG2	1:A:176:LEU:HD11	2.51	0.40
1:B:168:ALA:O	1:B:172:LEU:HG	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	272/275 (99%)	246 (90%)	22 (8%)	4 (2%)	13	22
1	B	272/275 (99%)	252 (93%)	16 (6%)	4 (2%)	13	22
1	C	272/275 (99%)	251 (92%)	16 (6%)	5 (2%)	11	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	272/275 (99%)	254 (93%)	15 (6%)	3 (1%)	17	31
All	All	1088/1100 (99%)	1003 (92%)	69 (6%)	16 (2%)	13	22

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	259	GLU
1	B	274	GLU
1	D	259	GLU
1	C	146	LEU
1	C	259	GLU
1	C	274	GLU
1	A	258	VAL
1	A	274	GLU
1	B	146	LEU
1	C	54	GLU
1	A	259	GLU
1	A	263	ASN
1	D	260	GLU
1	D	258	VAL
1	B	258	VAL
1	C	258	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	226/227 (100%)	208 (92%)	18 (8%)	15	28
1	B	226/227 (100%)	213 (94%)	13 (6%)	25	45
1	C	226/227 (100%)	210 (93%)	16 (7%)	18	34
1	D	226/227 (100%)	211 (93%)	15 (7%)	21	38
All	All	904/908 (100%)	842 (93%)	62 (7%)	19	35

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	20	ILE
1	A	42	GLU
1	A	49	ARG
1	A	56	ASN
1	A	72	LYS
1	A	91	SER
1	A	100	LEU
1	A	104	LEU
1	A	133	LEU
1	A	146	LEU
1	A	157	VAL
1	A	203	ARG
1	A	208	TRP
1	A	229	MET
1	A	231	LEU
1	A	235	LEU
1	A	275	GLN
1	B	16	ASN
1	B	49	ARG
1	B	67	LYS
1	B	72	LYS
1	B	85	LEU
1	B	104	LEU
1	B	146	LEU
1	B	203	ARG
1	B	208	TRP
1	B	229	MET
1	B	231	LEU
1	B	240	SER
1	B	275	GLN
1	C	16	ASN
1	C	46	LYS
1	C	49	ARG
1	C	67	LYS
1	C	72	LYS
1	C	85	LEU
1	C	104	LEU
1	C	133	LEU
1	C	146	LEU
1	C	151	TYR
1	C	203	ARG
1	C	208	TRP

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Mol	Chain	Res	Type
1	C	229	MET
1	C	231	LEU
1	C	240	SER
1	C	266	THR
1	D	16	ASN
1	D	46	LYS
1	D	49	ARG
1	D	56	ASN
1	D	58	PRO
1	D	63	LEU
1	D	67	LYS
1	D	72	LYS
1	D	85	LEU
1	D	104	LEU
1	D	133	LEU
1	D	146	LEU
1	D	208	TRP
1	D	229	MET
1	D	231	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	26	GLN
1	A	30	ASN
1	A	53	GLN
1	A	76	ASN
1	A	128	ASN
1	A	154	HIS
1	A	156	ASN
1	A	184	ASN
1	A	226	ASN
1	B	16	ASN
1	B	30	ASN
1	B	53	GLN
1	B	56	ASN
1	B	76	ASN
1	B	128	ASN
1	B	156	ASN
1	B	184	ASN
1	B	226	ASN

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Mol	Chain	Res	Type
1	C	16	ASN
1	C	30	ASN
1	C	53	GLN
1	C	56	ASN
1	C	76	ASN
1	C	156	ASN
1	C	180	HIS
1	C	184	ASN
1	C	209	ASN
1	C	226	ASN
1	D	16	ASN
1	D	30	ASN
1	D	53	GLN
1	D	56	ASN
1	D	76	ASN
1	D	128	ASN
1	D	156	ASN
1	D	179	HIS
1	D	184	ASN
1	D	218	ASN
1	D	226	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

8 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	NAD	A	1780	-	38,48,48	1.68	3 (7%)	47,73,73	2.38	8 (17%)
3	TCL	A	2414	-	18,18,18	2.12	4 (22%)	25,25,25	0.98	3 (12%)
2	NAD	B	2780	-	38,48,48	1.66	3 (7%)	47,73,73	2.27	7 (14%)
3	TCL	B	3414	-	18,18,18	1.96	3 (16%)	25,25,25	0.77	0
2	NAD	C	3780	-	38,48,48	1.66	3 (7%)	47,73,73	2.25	10 (21%)
3	TCL	C	4414	-	18,18,18	2.03	5 (27%)	25,25,25	0.83	1 (4%)
2	NAD	D	4780	-	38,48,48	1.67	3 (7%)	47,73,73	2.37	10 (21%)
3	TCL	D	5414	-	18,18,18	2.10	4 (22%)	25,25,25	0.88	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	A	1780	-	-	0/22/62/62	0/5/5/5
3	TCL	A	2414	-	-	0/4/4/4	0/2/2/2
2	NAD	B	2780	-	-	0/22/62/62	0/5/5/5
3	TCL	B	3414	-	-	0/4/4/4	0/2/2/2
2	NAD	C	3780	-	-	0/22/62/62	0/5/5/5
3	TCL	C	4414	-	-	0/4/4/4	0/2/2/2
2	NAD	D	4780	-	-	0/22/62/62	0/5/5/5
3	TCL	D	5414	-	-	0/4/4/4	0/2/2/2

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	4414	TCL	C2-CL14	2.05	1.79	1.74
3	C	4414	TCL	C9-CL16	2.10	1.78	1.73
3	D	5414	TCL	C11-CL15	2.18	1.79	1.74
3	C	4414	TCL	C11-CL15	2.21	1.79	1.74
3	B	3414	TCL	C9-CL16	2.27	1.79	1.73
3	A	2414	TCL	C9-CL16	2.35	1.79	1.73
3	A	2414	TCL	C11-CL15	2.36	1.79	1.74
2	D	4780	NAD	C2A-N1A	2.46	1.38	1.33
2	A	1780	NAD	C2A-N1A	2.52	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2780	NAD	C2A-N1A	2.56	1.38	1.33
2	C	3780	NAD	C2A-N1A	2.66	1.39	1.33
3	D	5414	TCL	C9-CL16	2.73	1.80	1.73
2	A	1780	NAD	C2A-N3A	3.43	1.38	1.32
2	B	2780	NAD	C2A-N3A	3.51	1.38	1.32
2	D	4780	NAD	C2A-N3A	3.81	1.38	1.32
2	C	3780	NAD	C2A-N3A	3.91	1.39	1.32
3	B	3414	TCL	C8-C9	4.89	1.48	1.39
3	A	2414	TCL	C8-C9	5.12	1.48	1.39
3	C	4414	TCL	C8-C9	5.14	1.48	1.39
3	C	4414	TCL	C6-C5	5.22	1.48	1.40
3	D	5414	TCL	C6-C5	5.22	1.48	1.40
3	D	5414	TCL	C8-C9	5.23	1.49	1.39
3	B	3414	TCL	C6-C5	5.31	1.48	1.40
3	A	2414	TCL	C6-C5	5.67	1.49	1.40
2	D	4780	NAD	O7N-C7N	7.59	1.40	1.24
2	C	3780	NAD	O7N-C7N	7.80	1.40	1.24
2	B	2780	NAD	O7N-C7N	7.86	1.40	1.24
2	A	1780	NAD	O7N-C7N	8.14	1.41	1.24

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4780	NAD	N3A-C2A-N1A	-12.62	119.23	128.89
2	A	1780	NAD	N3A-C2A-N1A	-12.28	119.49	128.89
2	B	2780	NAD	N3A-C2A-N1A	-12.23	119.53	128.89
2	C	3780	NAD	N3A-C2A-N1A	-11.29	120.25	128.89
2	C	3780	NAD	C4B-O4B-C1B	-3.46	105.92	109.72
2	B	2780	NAD	PN-O3-PA	-3.44	123.07	132.73
2	C	3780	NAD	PN-O3-PA	-2.91	124.57	132.73
2	A	1780	NAD	PN-O3-PA	-2.88	124.65	132.73
2	B	2780	NAD	C5D-C4D-C3D	-2.72	104.41	115.21
2	D	4780	NAD	O7N-C7N-N7N	-2.64	118.87	122.59
2	A	1780	NAD	C4B-O4B-C1B	-2.59	106.87	109.72
2	A	1780	NAD	C1B-N9A-C4A	-2.58	123.05	126.94
2	A	1780	NAD	C2B-C1B-N9A	-2.50	110.48	114.29
2	D	4780	NAD	PN-O3-PA	-2.47	125.78	132.73
2	D	4780	NAD	C4B-O4B-C1B	-2.45	107.03	109.72
2	D	4780	NAD	C5D-C4D-C3D	-2.41	105.65	115.21
2	C	3780	NAD	C2B-C1B-N9A	-2.20	110.93	114.29
2	A	1780	NAD	C5D-C4D-C3D	-2.16	106.64	115.21
2	D	4780	NAD	C4A-C5A-N7A	-2.08	107.56	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3780	NAD	C1B-N9A-C4A	-2.05	123.85	126.94
2	B	2780	NAD	C4B-O4B-C1B	-2.04	107.47	109.72
3	A	2414	TCL	O7-C5-C6	2.02	120.09	116.12
2	C	3780	NAD	O2A-PA-O3	2.07	114.46	105.09
3	C	4414	TCL	C13-C12-C11	2.21	121.68	119.23
3	A	2414	TCL	C13-C12-C11	2.30	121.79	119.23
2	D	4780	NAD	O3-PN-O5D	2.33	109.11	102.94
2	D	4780	NAD	C3N-C7N-N7N	2.42	120.47	117.82
3	A	2414	TCL	C8-O7-C5	2.43	123.66	117.75
2	B	2780	NAD	C3N-C7N-N7N	2.59	120.65	117.82
2	B	2780	NAD	O4B-C1B-N9A	2.66	113.67	108.10
2	C	3780	NAD	O3-PN-O5D	2.67	110.03	102.94
2	C	3780	NAD	C3N-C7N-N7N	2.81	120.89	117.82
2	A	1780	NAD	O4B-C1B-N9A	3.05	114.49	108.10
2	D	4780	NAD	O4B-C1B-N9A	3.16	114.70	108.10
2	C	3780	NAD	O4B-C1B-N9A	3.32	115.05	108.10
2	D	4780	NAD	O4D-C1D-N1N	4.27	112.82	108.13
2	C	3780	NAD	O4D-C1D-N1N	4.29	112.84	108.13
2	B	2780	NAD	O4D-C1D-N1N	4.31	112.87	108.13
2	A	1780	NAD	O4D-C1D-N1N	6.22	114.97	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1780	NAD	1	0
3	A	2414	TCL	2	0
2	B	2780	NAD	6	0
3	B	3414	TCL	2	0
2	C	3780	NAD	2	0
3	C	4414	TCL	3	0
2	D	4780	NAD	4	0
3	D	5414	TCL	2	0

## 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	274/275 (99%)	0.57	24 (8%) 12 13	17, 36, 64, 84	0
1	B	274/275 (99%)	0.44	24 (8%) 12 13	16, 30, 56, 77	0
1	C	274/275 (99%)	0.45	19 (6%) 20 22	15, 31, 54, 73	0
1	D	274/275 (99%)	0.51	21 (7%) 16 18	16, 33, 61, 80	0
All	All	1096/1100 (99%)	0.49	88 (8%) 15 16	15, 32, 61, 84	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	273	LYS	7.7
1	A	274	GLU	7.2
1	D	274	GLU	7.0
1	A	272	HIS	6.1
1	B	260	GLU	6.1
1	A	198	GLY	6.1
1	C	200	ALA	6.0
1	C	273	LYS	5.8
1	D	273	LYS	5.6
1	C	274	GLU	5.6
1	B	274	GLU	5.5
1	B	259	GLU	5.3
1	D	263	ASN	5.2
1	A	275	GLN	5.1
1	A	259	GLU	5.1
1	D	260	GLU	4.9
1	D	275	GLN	4.6
1	C	263	ASN	4.6
1	D	259	GLU	4.6
1	C	260	GLU	4.5
1	B	275	GLN	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	261	LYS	4.4
1	C	261	LYS	4.3
1	B	273	LYS	4.2
1	A	260	GLU	4.2
1	C	259	GLU	4.1
1	A	200	ALA	4.0
1	A	263	ASN	3.8
1	B	264	LYS	3.7
1	C	275	GLN	3.6
1	C	3	PHE	3.6
1	D	261	LYS	3.6
1	C	201	ASP	3.6
1	A	262	ASP	3.5
1	D	272	HIS	3.5
1	B	263	ASN	3.5
1	A	3	PHE	3.3
1	B	262	ASP	3.3
1	A	199	ILE	3.3
1	B	200	ALA	3.3
1	B	272	HIS	3.2
1	D	258	VAL	3.2
1	C	264	LYS	3.1
1	B	201	ASP	3.0
1	C	198	GLY	3.0
1	C	78	VAL	2.9
1	C	69	GLU	2.9
1	A	265	ALA	2.9
1	B	261	LYS	2.8
1	A	258	VAL	2.8
1	A	264	LYS	2.7
1	D	201	ASP	2.7
1	D	199	ILE	2.7
1	B	194	LEU	2.7
1	D	262	ASP	2.6
1	D	208	TRP	2.6
1	B	3	PHE	2.5
1	B	207	LYS	2.5
1	B	2	GLY	2.5
1	D	78	VAL	2.4
1	B	68	GLU	2.4
1	B	69	GLU	2.4
1	A	208	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	46	LYS	2.4
1	D	198	GLY	2.3
1	C	257	ALA	2.3
1	B	179	HIS	2.3
1	D	257	ALA	2.2
1	D	264	LYS	2.2
1	C	6	GLY	2.1
1	D	266	THR	2.1
1	D	80	LYS	2.1
1	C	262	ASP	2.1
1	D	10	LEU	2.1
1	A	197	SER	2.1
1	B	97	LYS	2.1
1	D	46	LYS	2.1
1	B	269	TRP	2.1
1	C	79	LYS	2.1
1	C	199	ILE	2.1
1	A	98	GLU	2.1
1	B	197	SER	2.1
1	B	180	HIS	2.1
1	A	257	ALA	2.0
1	A	46	LYS	2.0
1	A	80	LYS	2.0
1	A	76	ASN	2.0
1	A	269	TRP	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	TCL	A	2414	17/17	0.84	0.23	0.85	42,45,46,46	0
3	TCL	D	5414	17/17	0.88	0.19	0.46	29,33,34,34	0
2	NAD	C	3780	44/44	0.92	0.17	0.34	32,36,39,39	0
2	NAD	B	2780	44/44	0.94	0.16	0.03	30,34,36,36	0
3	TCL	B	3414	17/17	0.92	0.15	-0.10	30,34,35,35	0
2	NAD	D	4780	44/44	0.94	0.15	-0.38	31,41,44,44	0
2	NAD	A	1780	44/44	0.92	0.15	-0.51	36,39,44,45	0
3	TCL	C	4414	17/17	0.92	0.14	-0.61	38,40,41,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.