



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 02:01 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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

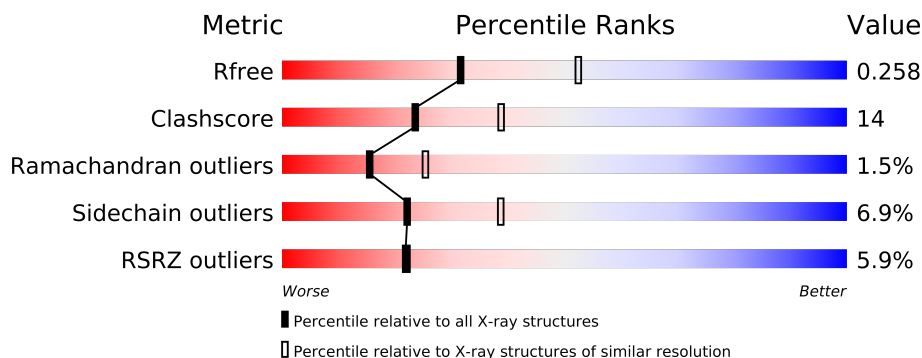
The following versions of software and data (see [references](#)) were used in the production of this report:

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

1 Overall quality at a glance

The reported resolution of this entry is 2.50 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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

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

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 8866 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called 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: $C_{21}H_{27}N_7O_{14}P_2$).



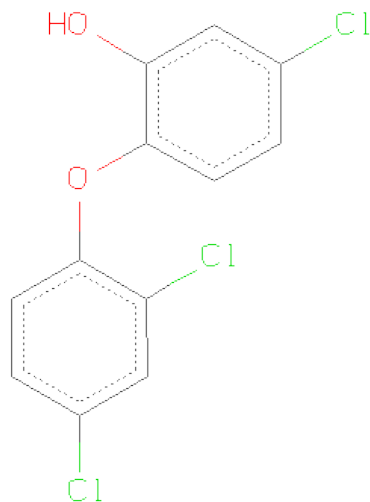
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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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: C₁₂H₇Cl₃O₂).



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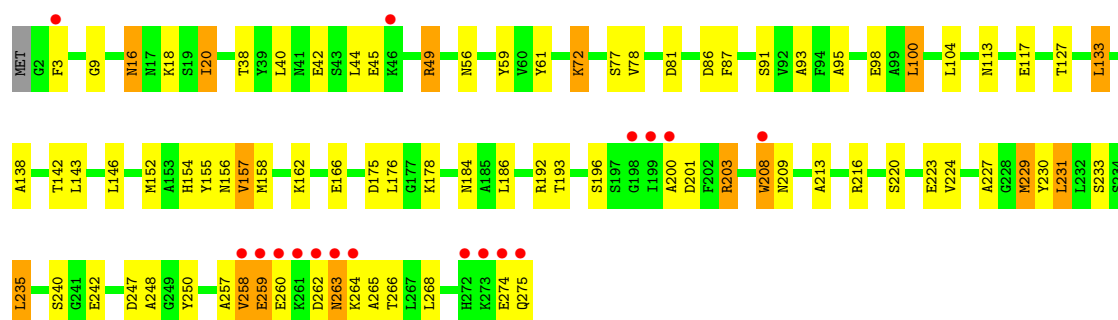
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	59	Total	O	0	0
			59	59		

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.

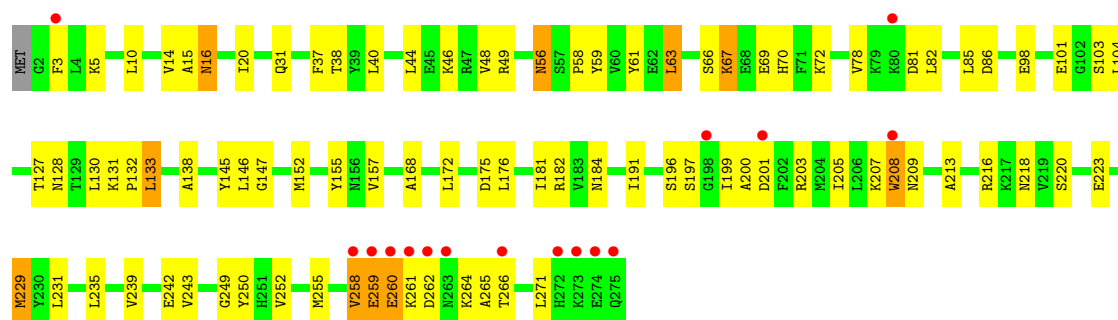
- Molecule 1: Enoyl-[acyl-carrier-protein]reductase [NADH]

Chain A: 



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

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	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 ²)	34.1	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 32.0	EDS
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	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 ²)	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.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2102	0	2125	65	0
1	B	2102	0	2125	61	0
1	C	2102	0	2125	65	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2102	0	2125	66	0
2	A	44	0	26	1	0
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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (249) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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	Distance(Å)	Clash(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
1:D:152:MET:CE	4:D:5416:HOH:O	2.51	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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

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	272/275 (99%)	246 (90%)	22 (8%)	4 (2%)	15	25
1	B	272/275 (99%)	252 (93%)	16 (6%)	4 (2%)	15	25
1	C	272/275 (99%)	251 (92%)	16 (6%)	5 (2%)	13	20
1	D	272/275 (99%)	254 (93%)	15 (6%)	3 (1%)	21	34
All	All	1088/1100 (99%)	1003 (92%)	69 (6%)	16 (2%)	15	25

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	259	GLU

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Mol	Chain	Res	Type
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%)	17	31
1	B	226/227 (100%)	213 (94%)	13 (6%)	28	49
1	C	226/227 (100%)	210 (93%)	16 (7%)	21	37
1	D	226/227 (100%)	211 (93%)	15 (7%)	24	41
All	All	904/908 (100%)	842 (93%)	62 (7%)	22	39

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

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Mol	Chain	Res	Type
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
1	C	229	MET
1	C	231	LEU
1	C	240	SER
1	C	266	THR
1	D	16	ASN
1	D	46	LYS

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Mol	Chain	Res	Type
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
1	C	16	ASN
1	C	30	ASN
1	C	53	GLN
1	C	56	ASN
1	C	76	ASN
1	C	156	ASN

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Mol	Chain	Res	Type
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 ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	-	48,48,48	1.40	3 (6%)	73,73,73	2.08	13 (17%)
3	TCL	A	2414	-	18,18,18	2.04	4 (22%)	25,25,25	0.95	2 (8%)
2	NAD	B	2780	-	48,48,48	1.42	4 (8%)	73,73,73	1.94	12 (16%)
3	TCL	B	3414	-	18,18,18	1.89	3 (16%)	25,25,25	0.74	0
2	NAD	C	3780	-	48,48,48	1.40	3 (6%)	73,73,73	1.99	10 (13%)
3	TCL	C	4414	-	18,18,18	1.96	5 (27%)	25,25,25	0.81	1 (4%)
2	NAD	D	4780	-	48,48,48	1.42	4 (8%)	73,73,73	2.07	13 (17%)
3	TCL	D	5414	-	18,18,18	2.02	4 (22%)	25,25,25	0.85	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	-	-	0/30/62/62	0/3/5/5
3	TCL	A	2414	-	-	0/4/4/4	0/2/2/2
2	NAD	B	2780	-	-	0/30/62/62	0/3/5/5
3	TCL	B	3414	-	-	0/4/4/4	0/2/2/2
2	NAD	C	3780	-	-	0/30/62/62	0/3/5/5
3	TCL	C	4414	-	-	0/4/4/4	0/2/2/2
2	NAD	D	4780	-	-	0/30/62/62	0/3/5/5
3	TCL	D	5414	-	-	0/4/4/4	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1780	NAD	O7N-C7N	7.40	1.41	1.24
2	B	2780	NAD	O7N-C7N	7.15	1.40	1.24
2	C	3780	NAD	O7N-C7N	7.10	1.40	1.24
2	D	4780	NAD	O7N-C7N	6.91	1.40	1.24
3	A	2414	TCL	C6-C5	5.50	1.49	1.40
3	B	3414	TCL	C6-C5	5.15	1.48	1.40
3	D	5414	TCL	C6-C5	5.06	1.48	1.40
3	C	4414	TCL	C6-C5	5.06	1.48	1.40
3	D	5414	TCL	C8-C9	4.94	1.49	1.39
3	C	4414	TCL	C8-C9	4.86	1.48	1.39
3	A	2414	TCL	C8-C9	4.84	1.48	1.39
3	B	3414	TCL	C8-C9	4.62	1.48	1.39
2	C	3780	NAD	C2A-N3A	3.51	1.39	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	4780	NAD	C2A-N3A	3.42	1.38	1.32
2	B	2780	NAD	C2A-N3A	3.15	1.38	1.32
2	A	1780	NAD	C2A-N3A	3.07	1.38	1.32
3	D	5414	TCL	C9-CL16	2.63	1.80	1.73
2	C	3780	NAD	C2A-N1A	2.56	1.39	1.33
2	B	2780	NAD	C2A-N1A	2.47	1.38	1.33
2	B	2780	NAD	C2N-N1N	2.46	1.38	1.35
2	A	1780	NAD	C2A-N1A	2.42	1.38	1.33
2	D	4780	NAD	C2A-N1A	2.37	1.38	1.33
3	A	2414	TCL	C11-CL15	2.31	1.79	1.74
3	A	2414	TCL	C9-CL16	2.26	1.79	1.73
3	B	3414	TCL	C9-CL16	2.19	1.79	1.73
3	C	4414	TCL	C11-CL15	2.16	1.79	1.74
2	D	4780	NAD	C2N-N1N	2.15	1.38	1.35
3	D	5414	TCL	C11-CL15	2.13	1.79	1.74
3	C	4414	TCL	C9-CL16	2.02	1.78	1.73
3	C	4414	TCL	C2-CL14	2.00	1.79	1.74

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4780	NAD	N3A-C2A-N1A	-11.33	119.23	128.71
2	A	1780	NAD	N3A-C2A-N1A	-11.02	119.49	128.71
2	B	2780	NAD	N3A-C2A-N1A	-10.98	119.53	128.71
2	C	3780	NAD	N3A-C2A-N1A	-10.12	120.25	128.71
2	C	3780	NAD	O4B-C1B-N9A	7.10	115.05	108.44
2	A	1780	NAD	O4D-C1D-N1N	6.86	114.97	107.95
2	D	4780	NAD	O4B-C1B-N9A	6.73	114.70	108.44
2	A	1780	NAD	O4B-C1B-N9A	6.50	114.49	108.44
2	B	2780	NAD	O4B-C1B-N9A	5.62	113.67	108.44
2	B	2780	NAD	O4D-C1D-N1N	4.81	112.87	107.95
2	C	3780	NAD	O4D-C1D-N1N	4.78	112.84	107.95
2	D	4780	NAD	O4D-C1D-N1N	4.76	112.82	107.95
2	C	3780	NAD	N3A-C4A-N9A	3.79	132.27	125.43
2	D	4780	NAD	N3A-C4A-N9A	3.59	131.91	125.43
2	A	1780	NAD	N3A-C4A-N9A	3.58	131.90	125.43
2	B	2780	NAD	N3A-C4A-N9A	3.56	131.85	125.43
2	C	3780	NAD	C4B-O4B-C1B	-3.53	105.92	109.75
2	C	3780	NAD	C3N-C7N-N7N	2.74	120.89	117.77
2	B	2780	NAD	C5D-C4D-C3D	-2.70	104.41	115.21
2	A	1780	NAD	C4B-O4B-C1B	-2.65	106.87	109.75
2	D	4780	NAD	O7N-C7N-N7N	-2.57	118.87	122.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2780	NAD	C3N-C7N-N7N	2.53	120.65	117.77
2	D	4780	NAD	C4B-O4B-C1B	-2.50	107.03	109.75
2	D	4780	NAD	C2A-N3A-C4A	2.45	120.98	114.01
2	D	4780	NAD	O4D-C1D-C2D	-2.43	103.05	106.77
2	D	4780	NAD	C5A-C4A-N3A	-2.42	120.43	125.70
2	A	1780	NAD	N7A-C8A-N9A	-2.42	107.53	114.36
2	D	4780	NAD	C5D-C4D-C3D	-2.39	105.65	115.21
2	A	1780	NAD	C8A-N9A-C4A	2.38	108.72	106.90
2	D	4780	NAD	C3N-C7N-N7N	2.37	120.47	117.77
3	A	2414	TCL	C8-O7-C5	2.33	123.66	117.93
3	A	2414	TCL	C13-C12-C11	2.30	121.79	119.22
2	B	2780	NAD	PN-O3-PA	-2.30	123.07	132.95
2	D	4780	NAD	C4A-C5A-N7A	-2.28	107.56	109.52
2	C	3780	NAD	C5A-C4A-N3A	-2.24	120.83	125.70
2	C	3780	NAD	N7A-C8A-N9A	-2.22	108.09	114.36
2	B	2780	NAD	C2A-N3A-C4A	2.21	120.31	114.01
3	C	4414	TCL	C13-C12-C11	2.21	121.68	119.22
2	B	2780	NAD	C5A-C4A-N3A	-2.17	120.97	125.70
2	C	3780	NAD	O3-PN-O5D	2.16	110.03	101.36
2	A	1780	NAD	C2A-N3A-C4A	2.16	120.16	114.01
2	A	1780	NAD	C4A-C5A-N7A	-2.15	107.68	109.52
2	A	1780	NAD	C5D-C4D-C3D	-2.14	106.64	115.21
2	B	2780	NAD	N7A-C8A-N9A	-2.12	108.37	114.36
2	B	2780	NAD	C4B-O4B-C1B	-2.09	107.47	109.75
2	A	1780	NAD	C6N-N1N-C2N	-2.09	119.67	122.04
2	A	1780	NAD	C1B-N9A-C4A	-2.07	123.05	126.64
2	C	3780	NAD	C2A-N3A-C4A	2.05	119.85	114.01
2	B	2780	NAD	O4D-C1D-C2D	-2.05	103.63	106.77
2	A	1780	NAD	C5A-C4A-N3A	-2.04	121.25	125.70
2	D	4780	NAD	N7A-C8A-N9A	-2.03	108.62	114.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	274/275 (99%)	0.49	17 (6%) 20 20	17, 36, 64, 84	0
1	B	274/275 (99%)	0.35	17 (6%) 20 20	16, 30, 56, 77	0
1	C	274/275 (99%)	0.36	15 (5%) 24 24	15, 31, 54, 73	0
1	D	274/275 (99%)	0.40	16 (5%) 22 23	16, 33, 61, 80	0
All	All	1096/1100 (99%)	0.40	65 (5%) 22 22	15, 32, 61, 84	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	273	LYS	7.5
1	A	274	GLU	7.0
1	D	274	GLU	6.7
1	D	273	LYS	5.6
1	C	273	LYS	5.6
1	C	200	ALA	5.3
1	C	274	GLU	5.2
1	B	259	GLU	5.0
1	B	274	GLU	4.9
1	B	260	GLU	4.9
1	A	272	HIS	4.8
1	A	275	GLN	4.7
1	D	260	GLU	4.7
1	A	261	LYS	4.5
1	B	275	GLN	4.4
1	A	198	GLY	4.4
1	A	259	GLU	4.4
1	B	273	LYS	4.4
1	D	263	ASN	4.3
1	D	259	GLU	4.3
1	C	261	LYS	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	259	GLU	4.1
1	C	263	ASN	4.1
1	D	261	LYS	4.0
1	D	275	GLN	4.0
1	A	260	GLU	3.8
1	C	260	GLU	3.8
1	C	275	GLN	3.8
1	A	199	ILE	3.7
1	A	200	ALA	3.7
1	C	3	PHE	3.6
1	A	263	ASN	3.5
1	A	3	PHE	3.3
1	B	264	LYS	3.3
1	B	263	ASN	3.1
1	A	262	ASP	3.0
1	C	264	LYS	3.0
1	B	262	ASP	3.0
1	B	261	LYS	3.0
1	D	258	VAL	2.9
1	B	200	ALA	2.8
1	C	201	ASP	2.7
1	B	3	PHE	2.7
1	D	208	TRP	2.6
1	D	272	HIS	2.6
1	B	201	ASP	2.6
1	A	264	LYS	2.6
1	B	272	HIS	2.5
1	B	46	LYS	2.3
1	A	258	VAL	2.3
1	B	194	LEU	2.3
1	C	69	GLU	2.3
1	C	207	LYS	2.2
1	B	207	LYS	2.2
1	B	2	GLY	2.2
1	D	3	PHE	2.2
1	D	201	ASP	2.2
1	A	208	TRP	2.2
1	D	262	ASP	2.2
1	C	198	GLY	2.1
1	D	266	THR	2.1
1	C	208	TRP	2.0
1	A	46	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	80	LYS	2.0
1	D	198	GLY	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	TCL	D	5414	17/17	0.20	0.92	29,33,34,34	0
3	TCL	A	2414	17/17	0.23	0.64	42,45,46,46	0
2	NAD	C	3780	44/44	0.16	0.21	32,36,39,39	0
2	NAD	B	2780	44/44	0.15	0.02	30,34,36,36	0
3	TCL	B	3414	17/17	0.14	-0.01	30,34,35,35	0
2	NAD	D	4780	44/44	0.14	-0.42	31,41,44,44	0
3	TCL	C	4414	17/17	0.14	-0.60	38,40,41,43	0
2	NAD	A	1780	44/44	0.15	-0.61	36,39,44,45	0

6.5 Other polymers

There are no such residues in this entry.