



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:22 PM GMT

PDB ID : 4RQU
Title : Alcohol Dehydrogenase crystal structure in complex with NAD
Authors : Xu, Y.W.
Deposited on : 2014-11-05
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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

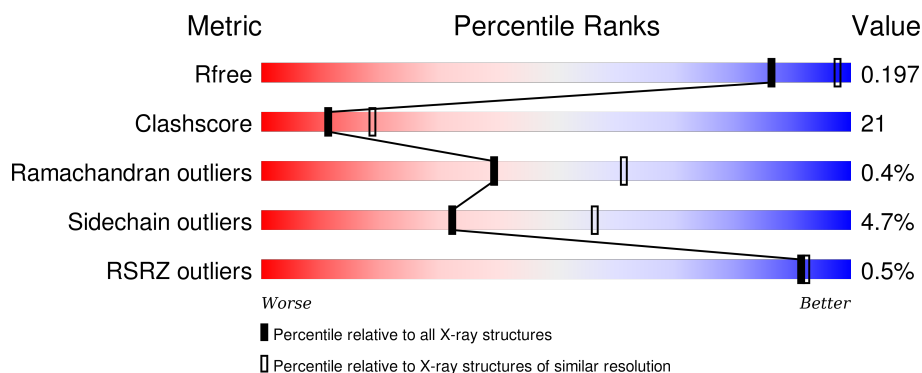
1 Overall quality at a glance

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 ≥ 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	375	 71% 26% ..
1	B	375	 % 55% 38% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11545 atoms, of which 5700 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 Alcohol Dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	375	Total	C	H	N	O	S	0	0	0
			5700	1812	2837	491	540	20			
1	B	375	Total	C	H	N	O	S	0	0	0
			5700	1812	2837	491	540	20			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	SER	-	EXPRESSION TAG	UNP P06525
B	5	SER	-	EXPRESSION TAG	UNP P06525

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

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

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	P	0	0
			70	21	26	7	14	2		

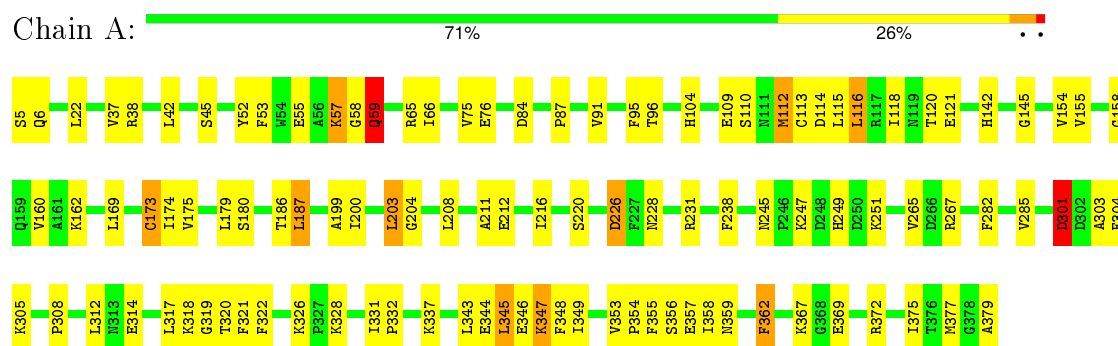
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	36	Total	O	0	0
			36	36		

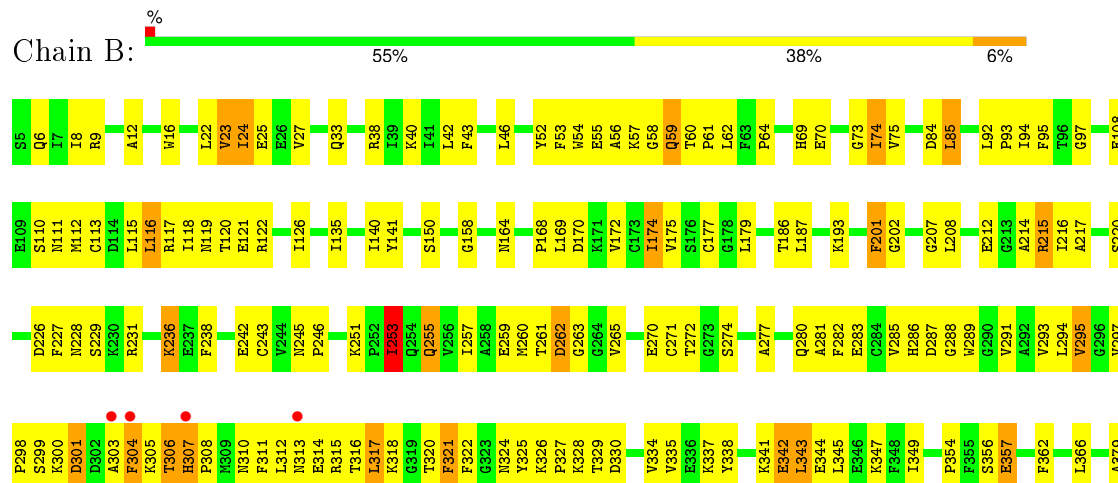
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: Alcohol Dehydrogenase



• Molecule 1: Alcohol Dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.15Å 101.15Å 165.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.37 – 2.50 48.37 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.4 (48.37-2.50) 96.3 (48.37-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.170 , 0.198 0.171 , 0.197	Depositor DCC
R_{free} test set	2388 reflections (7.72%)	DCC
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 20.9	EDS
Estimated twinning fraction	0.467 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtriage
Outliers	0 of 33347 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11545	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.92	1/2925 (0.0%)	1.11	15/3953 (0.4%)
1	B	1.01	3/2925 (0.1%)	1.21	19/3953 (0.5%)
All	All	0.97	4/5850 (0.1%)	1.16	34/7906 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	4
All	All	0	5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	321	PHE	CB-CG	-8.17	1.37	1.51
1	A	173	CYS	CB-SG	-5.44	1.73	1.81
1	B	113	CYS	CB-SG	-5.29	1.73	1.81
1	B	217	ALA	CA-CB	-5.02	1.42	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	215	ARG	NE-CZ-NH2	-12.19	114.20	120.30
1	B	201	PHE	CB-CG-CD1	-8.47	114.87	120.80
1	B	85	LEU	CB-CG-CD2	-8.41	96.70	111.00
1	B	253	ILE	CG1-CB-CG2	-7.42	95.07	111.40
1	A	174	ILE	CG1-CB-CG2	-7.41	95.09	111.40

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	201	PHE	CB-CG-CD2	7.40	125.98	120.80
1	B	300	LYS	CD-CE-NZ	7.36	128.63	111.70
1	B	215	ARG	CD-NE-CZ	6.86	133.21	123.60
1	A	66	ILE	CG1-CB-CG2	-6.79	96.46	111.40
1	A	58	GLY	N-CA-C	6.38	129.06	113.10
1	B	295	VAL	CG1-CB-CG2	-6.36	100.72	110.90
1	A	114	ASP	CB-CG-OD1	-6.30	112.63	118.30
1	A	345	LEU	CA-CB-CG	6.21	129.57	115.30
1	A	187	LEU	CB-CG-CD2	6.10	121.37	111.00
1	B	260	MET	CG-SD-CE	-5.84	90.86	100.20
1	B	297	VAL	CG1-CB-CG2	5.83	120.22	110.90
1	B	304	PHE	CB-CG-CD1	-5.81	116.73	120.80
1	B	301	ASP	CB-CG-OD1	-5.78	113.10	118.30
1	A	301	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	348	PHE	CB-CG-CD1	-5.76	116.77	120.80
1	B	215	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	372	ARG	NE-CZ-NH2	-5.65	117.47	120.30
1	B	208	LEU	CB-CG-CD2	-5.63	101.43	111.00
1	B	174	ILE	CG1-CB-CG2	-5.57	99.15	111.40
1	A	226	ASP	CB-CG-OD1	5.51	123.26	118.30
1	A	353	VAL	CG1-CB-CG2	-5.50	102.10	110.90
1	A	362	PHE	CB-CG-CD2	-5.45	116.98	120.80
1	B	366	LEU	CB-CG-CD2	5.33	120.07	111.00
1	B	9	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	347	LYS	CD-CE-NZ	5.25	123.78	111.70
1	A	59	GLN	N-CA-C	5.25	125.17	111.00
1	A	203	LEU	CB-CG-CD2	5.13	119.72	111.00
1	B	92	LEU	CB-CG-CD2	-5.10	102.32	111.00
1	B	46	LEU	CB-CG-CD1	5.06	119.61	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	55	GLU	Sidechain
1	B	164	ASN	Sidechain
1	B	271	CYS	Mainchain
1	B	33	GLN	Sidechain
1	B	357	GLU	Sidechain

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	2863	2837	2826	95	1
1	B	2863	2837	2825	162	1
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	B	44	26	24	0	0
4	A	35	0	0	13	0
4	B	36	0	0	11	0
All	All	5845	5700	5675	236	1

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 21.

All (236) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:319:GLY:HA3	1:B:311:PHE:HB3	1.41	1.01
1:B:280:GLN:NE2	4:B:532:HOH:O	1.92	1.00
1:B:59:GLN:NE2	1:B:121:GLU:OE2	2.02	0.91
1:A:249:HIS:ND1	4:A:533:HOH:O	2.02	0.91
1:A:308:PRO:HB2	1:B:298:PRO:HG3	1.49	0.91
1:A:251:LYS:O	4:A:512:HOH:O	1.89	0.90
1:A:228:ASN:O	4:A:523:HOH:O	1.89	0.90
1:B:174:ILE:HD12	1:B:179:LEU:HD12	1.53	0.89
1:A:226:ASP:OD1	4:A:523:HOH:O	1.90	0.88
1:B:320:THR:N	4:B:533:HOH:O	2.08	0.84
1:B:255:GLN:NE2	1:B:259:GLU:OE2	2.11	0.84
1:B:229:SER:OG	1:B:245:ASN:ND2	2.12	0.82
1:B:6:GLN:O	1:B:38:ARG:NH2	2.12	0.82
1:B:119:ASN:O	4:B:517:HOH:O	1.96	0.82
1:B:227:PHE:HE1	1:B:246:PRO:HB2	1.46	0.81
1:A:308:PRO:CB	1:B:298:PRO:HG3	2.10	0.80
1:B:342:GLU:O	4:B:516:HOH:O	2.02	0.77
1:B:287:ASP:HB3	1:B:315:ARG:CZ	2.14	0.77
1:A:305:LYS:HA	1:B:303:ALA:HB1	1.67	0.77
1:B:59:GLN:HE22	1:B:121:GLU:CD	1.88	0.76

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:TYR:OH	1:B:58:GLY:HA3	1.86	0.75
1:B:42:LEU:HD11	1:B:74:ILE:HG13	1.68	0.75
1:A:344:GLU:HB2	1:A:347:LYS:HE2	1.66	0.74
1:B:343:LEU:HD13	1:B:345:LEU:HD21	1.68	0.74
1:B:23:VAL:HG12	1:B:25:GLU:HG2	1.69	0.73
1:A:245:ASN:OD1	1:A:247:LYS:N	2.22	0.72
1:B:62:LEU:HD11	1:B:141:TYR:CD2	2.24	0.72
1:A:173:CYS:SG	4:A:507:HOH:O	2.48	0.72
1:B:174:ILE:HD12	1:B:179:LEU:CD1	2.19	0.71
1:B:349:ILE:O	4:B:508:HOH:O	2.10	0.70
1:B:270:GLU:O	1:B:295:VAL:HG12	1.91	0.70
1:B:287:ASP:HB3	1:B:315:ARG:NH2	2.07	0.69
1:B:56:ALA:HB1	1:B:61:PRO:HB3	1.74	0.69
1:B:282:PHE:HE2	1:B:306:THR:HG21	1.58	0.69
1:B:344:GLU:O	1:B:347:LYS:HE2	1.94	0.68
1:B:84:ASP:OD2	4:B:535:HOH:O	2.11	0.68
1:B:118:ILE:HB	1:B:321:PHE:HZ	1.59	0.68
1:B:27:VAL:HG12	1:B:135:ILE:CG1	2.24	0.68
1:A:95:PHE:HD1	1:A:118:ILE:HD12	1.61	0.66
1:B:311:PHE:O	1:B:314:GLU:N	2.26	0.66
1:A:320:THR:O	1:B:312:LEU:HD23	1.96	0.66
1:A:112:MET:SD	1:B:313:ASN:HA	2.36	0.65
1:B:316:THR:HG22	1:B:317:LEU:N	2.11	0.65
1:A:322:PHE:CE2	1:A:331:ILE:HD11	2.31	0.65
1:A:282:PHE:O	1:A:285:VAL:HG22	1.97	0.65
1:B:52:TYR:CE1	1:B:299:SER:HB3	2.32	0.64
1:B:261:THR:O	1:B:263:GLY:N	2.31	0.64
1:A:95:PHE:CD1	1:A:118:ILE:HD12	2.32	0.64
1:A:104:HIS:NE2	1:B:287:ASP:OD2	2.23	0.63
1:B:122:ARG:CZ	1:B:126:ILE:HD11	2.29	0.63
1:B:356:SER:OG	1:B:379:ALA:O	2.15	0.62
1:B:285:VAL:HG23	1:B:315:ARG:HD3	1.81	0.62
1:B:253:ILE:HD12	1:B:253:ILE:O	2.00	0.62
1:A:265:VAL:HG23	1:A:285:VAL:HG12	1.82	0.61
1:A:312:LEU:HD12	1:A:312:LEU:H	1.66	0.61
1:A:344:GLU:HB2	1:A:347:LYS:CE	2.31	0.61
1:B:288:GLY:N	1:B:313:ASN:O	2.32	0.61
1:B:27:VAL:HG12	1:B:135:ILE:HG13	1.83	0.61
1:A:145:GLY:O	1:A:155:VAL:HG13	2.01	0.61
1:B:291:VAL:HG22	1:B:316:THR:HB	1.83	0.61
1:A:305:LYS:CA	1:B:303:ALA:HB1	2.30	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:GLY:HA2	1:B:328:LYS:HD2	1.81	0.60
1:A:42:LEU:HD12	1:A:169:LEU:CD1	2.31	0.60
1:A:344:GLU:CB	1:A:347:LYS:HE2	2.31	0.60
1:B:282:PHE:CD2	1:B:304:PHE:HZ	2.20	0.60
1:A:367:LYS:HB2	1:A:369:GLU:HG3	1.84	0.60
1:A:303:ALA:HB1	1:B:305:LYS:HB3	1.84	0.59
1:B:54:TRP:CE3	1:B:55:GLU:HG3	2.36	0.59
1:B:55:GLU:OE1	1:B:57:LYS:HD2	2.04	0.58
1:A:179:LEU:HD11	1:A:345:LEU:CD2	2.33	0.58
1:B:62:LEU:HD11	1:B:141:TYR:CG	2.38	0.57
1:B:288:GLY:HA2	1:B:314:GLU:HB2	1.87	0.57
1:B:287:ASP:HA	1:B:315:ARG:CG	2.35	0.57
1:B:59:GLN:HE22	1:B:121:GLU:CG	2.17	0.57
1:A:305:LYS:HB3	1:B:303:ALA:CB	2.35	0.57
1:B:212:GLU:CD	1:B:215:ARG:HH21	2.07	0.57
1:A:312:LEU:HD12	1:A:312:LEU:N	2.19	0.56
1:A:226:ASP:N	4:A:515:HOH:O	2.37	0.56
1:A:245:ASN:O	4:A:533:HOH:O	2.18	0.56
1:A:42:LEU:HD12	1:A:169:LEU:HD12	1.87	0.56
1:B:320:THR:OG1	1:B:321:PHE:N	2.36	0.56
1:B:283:GLU:OE2	1:B:306:THR:OG1	2.06	0.56
1:A:304:PHE:O	1:B:305:LYS:HA	2.06	0.56
1:A:318:LYS:HG2	1:B:316:THR:HG23	1.88	0.55
1:B:43:PHE:CZ	1:B:170:ASP:HB3	2.42	0.55
1:B:261:THR:O	1:B:262:ASP:C	2.45	0.55
1:B:316:THR:CG2	1:B:317:LEU:N	2.69	0.55
1:B:115:LEU:HG	1:B:116:LEU:HD13	1.88	0.55
1:B:118:ILE:HB	1:B:321:PHE:CZ	2.41	0.55
1:B:110:SER:O	1:B:326:LYS:NZ	2.39	0.55
1:B:43:PHE:CE1	1:B:170:ASP:HB3	2.41	0.55
1:B:59:GLN:HE22	1:B:121:GLU:HG2	1.73	0.54
1:B:304:PHE:HB2	4:B:518:HOH:O	2.07	0.54
1:B:227:PHE:CE1	1:B:246:PRO:HB2	2.35	0.54
1:B:112:MET:SD	1:B:321:PHE:CD2	3.00	0.54
1:B:270:GLU:HB3	1:B:294:LEU:HA	1.90	0.54
1:A:346:GLU:HA	1:A:349:ILE:HD12	1.89	0.54
1:B:334:VAL:O	1:B:337:LYS:HB2	2.08	0.54
1:A:267:ARG:NH1	4:A:505:HOH:O	2.40	0.54
1:B:24:ILE:N	1:B:24:ILE:HD13	2.22	0.53
1:A:349:ILE:HG22	4:A:513:HOH:O	2.08	0.53
1:A:354:PRO:HG2	1:A:357:GLU:HG2	1.90	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:LYS:HE3	1:B:243:CYS:SG	2.49	0.52
1:B:304:PHE:N	4:B:518:HOH:O	2.41	0.52
1:B:55:GLU:CB	1:B:57:LYS:HG3	2.39	0.52
1:A:154:VAL:O	4:A:508:HOH:O	2.19	0.52
1:A:304:PHE:O	1:B:304:PHE:O	2.28	0.52
1:B:16:TRP:CE3	1:B:64:PRO:HB3	2.45	0.52
1:B:322:PHE:CG	1:B:322:PHE:O	2.63	0.52
1:B:212:GLU:OE1	1:B:347:LYS:NZ	2.42	0.52
1:B:285:VAL:O	1:B:315:ARG:NE	2.40	0.51
1:B:287:ASP:HB2	1:B:313:ASN:HB3	1.92	0.51
1:A:245:ASN:OD1	1:A:245:ASN:C	2.49	0.51
1:B:212:GLU:CD	1:B:215:ARG:NH2	2.64	0.51
1:B:357:GLU:HG3	4:B:534:HOH:O	2.11	0.51
1:A:179:LEU:HD11	1:A:345:LEU:HD23	1.92	0.51
1:A:91:VAL:HG12	1:A:162:LYS:HA	1.93	0.51
1:A:22:LEU:HD11	1:A:362:PHE:HZ	1.76	0.51
1:B:282:PHE:CE2	1:B:306:THR:HG21	2.42	0.51
1:A:321:PHE:HB2	1:B:312:LEU:O	2.11	0.51
1:B:193:LYS:NZ	4:B:505:HOH:O	2.38	0.51
1:A:158:GLY:HA2	1:A:328:LYS:HD2	1.92	0.50
1:B:53:PHE:CE1	1:B:120:THR:HB	2.47	0.50
1:B:255:GLN:O	1:B:259:GLU:HG3	2.10	0.50
1:B:238:PHE:CE1	1:B:347:LYS:HD2	2.47	0.50
1:B:287:ASP:CA	1:B:315:ARG:CZ	2.90	0.50
1:A:308:PRO:HB2	1:B:298:PRO:CG	2.32	0.49
1:B:97:GLY:HA3	1:B:116:LEU:HB3	1.93	0.49
1:B:42:LEU:HD11	1:B:74:ILE:CG1	2.39	0.49
1:B:201:PHE:N	1:B:201:PHE:CD1	2.80	0.49
1:B:212:GLU:O	1:B:216:ILE:HG13	2.13	0.49
1:B:55:GLU:HB2	1:B:57:LYS:HG3	1.94	0.49
1:A:113:CYS:SG	1:A:115:LEU:HB3	2.53	0.49
1:B:202:GLY:HA3	1:B:272:THR:HG22	1.95	0.48
1:B:69:HIS:CD2	1:B:177:CYS:SG	3.06	0.48
1:A:53:PHE:CZ	1:A:120:THR:HG21	2.48	0.48
1:B:27:VAL:HG12	1:B:135:ILE:HD11	1.94	0.48
1:B:228:ASN:O	1:B:231:ARG:HG3	2.12	0.48
1:B:287:ASP:CB	1:B:315:ARG:CZ	2.87	0.48
1:A:312:LEU:C	1:A:314:GLU:H	2.17	0.48
1:B:329:THR:HG22	1:B:330:ASP:OD1	2.14	0.48
1:A:45:SER:OG	4:A:507:HOH:O	2.20	0.48
1:A:200:ILE:CD1	1:A:211:ALA:HA	2.43	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:LYS:N	1:B:236:LYS:CD	2.77	0.48
1:B:305:LYS:O	1:B:306:THR:HG22	2.14	0.47
1:A:208:LEU:HD22	1:A:238:PHE:CD2	2.49	0.47
1:A:312:LEU:HA	4:B:533:HOH:O	2.14	0.47
1:A:317:LEU:HB3	1:B:317:LEU:HD23	1.95	0.47
1:B:40:LYS:O	1:B:74:ILE:N	2.40	0.47
1:A:75:VAL:HG12	1:A:87:PRO:HA	1.96	0.47
1:A:109:GLU:HB2	1:B:289:TRP:HB3	1.96	0.47
1:A:115:LEU:O	4:A:517:HOH:O	2.20	0.47
1:B:293:VAL:HA	1:B:318:LYS:O	2.14	0.47
1:B:112:MET:O	1:B:117:ARG:NH2	2.39	0.47
1:A:22:LEU:HD11	1:A:362:PHE:CZ	2.49	0.47
1:A:65:ARG:NE	1:A:142:HIS:O	2.35	0.47
1:B:343:LEU:HD13	1:B:345:LEU:CD2	2.40	0.46
1:A:59:GLN:OE1	1:A:121:GLU:HG2	2.16	0.46
1:B:70:GLU:HG2	1:B:177:CYS:HB3	1.98	0.46
1:A:343:LEU:HD21	1:A:345:LEU:HD21	1.98	0.46
1:A:204:GLY:HA2	1:A:231:ARG:NH1	2.31	0.46
1:B:93:PRO:O	1:B:94:ILE:HD13	2.16	0.46
1:B:169:LEU:HA	1:B:172:VAL:HG22	1.97	0.46
1:B:228:ASN:HB3	1:B:231:ARG:HG3	1.97	0.46
1:A:6:GLN:O	1:A:38:ARG:NH2	2.24	0.46
1:B:305:LYS:O	1:B:306:THR:CB	2.63	0.46
1:B:179:LEU:HD11	1:B:345:LEU:CD2	2.45	0.45
1:B:282:PHE:CD2	1:B:304:PHE:CZ	3.02	0.45
1:A:367:LYS:CB	1:A:369:GLU:HG3	2.47	0.45
1:B:24:ILE:N	1:B:24:ILE:CD1	2.78	0.45
1:B:179:LEU:HD11	1:B:345:LEU:HD22	1.99	0.45
1:B:27:VAL:HG12	1:B:135:ILE:CD1	2.46	0.45
1:A:38:ARG:HD2	1:A:76:GLU:OE1	2.16	0.45
1:B:22:LEU:HD11	1:B:362:PHE:CE2	2.53	0.45
1:B:52:TYR:CZ	1:B:58:GLY:HA3	2.52	0.44
1:B:338:TYR:CD1	1:B:338:TYR:O	2.70	0.44
1:B:179:LEU:CD1	1:B:345:LEU:HD22	2.48	0.44
1:B:57:LYS:NZ	1:B:57:LYS:HB3	2.32	0.44
1:B:201:PHE:O	1:B:272:THR:HG23	2.16	0.44
1:A:52:TYR:CE1	1:A:57:LYS:HG2	2.52	0.44
1:B:135:ILE:HD12	1:B:140:ILE:CD1	2.47	0.44
1:B:305:LYS:O	1:B:306:THR:HB	2.18	0.44
1:A:199:ALA:HB2	1:A:265:VAL:HG11	1.99	0.44
1:A:112:MET:HE2	1:A:112:MET:HB2	1.83	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ASP:OD1	1:A:84:ASP:N	2.42	0.43
1:B:325:TYR:O	1:B:327:PRO:HD3	2.18	0.43
1:A:175:VAL:O	1:A:180:SER:OG	2.27	0.43
1:B:226:ASP:OD1	1:B:227:PHE:N	2.51	0.43
1:B:23:VAL:HG11	1:B:25:GLU:OE2	2.18	0.43
1:A:53:PHE:CE1	1:A:120:THR:HB	2.54	0.43
1:B:186:THR:HG22	1:B:187:LEU:HD23	2.00	0.43
1:B:311:PHE:C	1:B:313:ASN:N	2.71	0.43
1:A:186:THR:HG22	1:A:187:LEU:HD23	2.00	0.43
1:B:111:ASN:HA	1:B:326:LYS:HE3	2.00	0.43
1:A:355:PHE:CD1	1:A:377:MET:SD	3.12	0.43
1:B:56:ALA:CB	1:B:61:PRO:HB3	2.44	0.43
1:A:337:LYS:HD3	1:A:337:LYS:HA	1.77	0.43
1:B:308:PRO:O	1:B:311:PHE:HB2	2.19	0.42
1:B:8:ILE:HD12	1:B:38:ARG:HD3	2.01	0.42
1:A:322:PHE:HE2	1:A:331:ILE:HD11	1.82	0.42
1:B:12:ALA:HA	1:B:150:SER:HA	2.01	0.42
1:A:203:LEU:HG	1:A:226:ASP:HB2	2.00	0.42
1:A:110:SER:O	1:A:326:LYS:NZ	2.45	0.42
1:B:175:VAL:HG12	1:B:175:VAL:O	2.19	0.42
1:B:286:HIS:C	1:B:315:ARG:HE	2.22	0.42
1:A:308:PRO:O	1:A:312:LEU:HD12	2.20	0.42
1:B:283:GLU:OE1	1:B:310:ASN:ND2	2.52	0.42
1:B:287:ASP:HA	1:B:315:ARG:HG3	2.02	0.42
1:A:331:ILE:N	1:A:332:PRO:CD	2.82	0.42
1:A:115:LEU:HG	1:A:116:LEU:HD13	2.00	0.42
1:B:277:ALA:O	1:B:281:ALA:N	2.48	0.42
1:A:319:GLY:CA	1:B:311:PHE:HB3	2.29	0.42
1:A:52:TYR:HD2	1:A:53:PHE:CD2	2.37	0.42
1:B:75:VAL:HG21	1:B:85:LEU:HD22	2.02	0.42
1:B:202:GLY:O	1:B:207:GLY:HA3	2.20	0.41
1:B:186:THR:HG21	1:B:214:ALA:N	2.35	0.41
1:A:42:LEU:HD12	1:A:169:LEU:HD13	2.02	0.41
1:A:96:THR:HG21	1:A:321:PHE:CZ	2.56	0.41
1:B:354:PRO:HG2	1:B:357:GLU:HG2	2.02	0.41
1:A:112:MET:SD	1:B:313:ASN:CA	3.06	0.41
1:B:70:GLU:CG	1:B:177:CYS:HB3	2.50	0.41
1:A:65:ARG:NH2	4:A:534:HOH:O	2.53	0.41
1:A:358:ILE:HG23	1:A:359:ASN:N	2.36	0.41
1:B:95:PHE:HE1	1:B:321:PHE:CE1	2.38	0.41
1:A:301:ASP:CG	1:B:307:HIS:HE2	2.23	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:GLU:O	1:A:216:ILE:HG13	2.21	0.41
1:A:37:VAL:HG21	1:A:160:VAL:HG21	2.02	0.41
1:A:305:LYS:HB3	1:B:303:ALA:HB2	2.02	0.41
1:B:55:GLU:HB3	1:B:57:LYS:HG3	2.02	0.41
1:B:42:LEU:HG	1:B:73:GLY:HA2	2.02	0.40
1:B:227:PHE:CE1	1:B:246:PRO:HG2	2.56	0.40
1:B:261:THR:HG21	1:B:265:VAL:HG13	2.03	0.40
1:A:355:PHE:HA	1:A:375:ILE:CG2	2.51	0.40
1:B:168:PRO:HB2	1:B:170:ASP:OD1	2.21	0.40
1:B:338:TYR:CD1	1:B:338:TYR:C	2.94	0.40
1:A:356:SER:HB3	1:A:379:ALA:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:SER:O	1:B:141:TYR:OH[6_554]	2.17	0.03

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	373/375 (100%)	359 (96%)	13 (4%)	1 (0%)	46	68
1	B	373/375 (100%)	357 (96%)	14 (4%)	2 (0%)	34	55
All	All	746/750 (100%)	716 (96%)	27 (4%)	3 (0%)	39	61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	262	ASP
1	B	306	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	59	GLN

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	310/310 (100%)	304 (98%)	6 (2%)	65	87
1	B	310/310 (100%)	287 (93%)	23 (7%)	17	31
All	All	620/620 (100%)	591 (95%)	29 (5%)	32	56

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	SER
1	A	57	LYS
1	A	59	GLN
1	A	112	MET
1	A	116	LEU
1	A	301	ASP
1	B	23	VAL
1	B	24	ILE
1	B	59	GLN
1	B	60	THR
1	B	74	ILE
1	B	108	GLU
1	B	116	LEU
1	B	220	SER
1	B	236	LYS
1	B	242	GLU
1	B	251	LYS
1	B	253	ILE
1	B	255	GLN
1	B	257	ILE
1	B	274	SER
1	B	301	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	307	HIS
1	B	317	LEU
1	B	324	ASN
1	B	335	VAL
1	B	341	LYS
1	B	342	GLU
1	B	343	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	35	HIS
1	B	59	GLN
1	B	245	ASN
1	B	255	GLN
1	B	324	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](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAD	B	403	-	38,48,48	3.63	12 (31%)	47,73,73	2.68	16 (34%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAD	B	403	-	-	0/22/62/62	0/5/5/5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	NAD	O4B-C4B	-6.40	1.30	1.45
3	B	403	NAD	O4D-C4D	-5.45	1.32	1.45
3	B	403	NAD	O3D-C3D	-3.94	1.33	1.43
3	B	403	NAD	O3B-C3B	-3.36	1.34	1.43
3	B	403	NAD	C5A-C4A	-2.21	1.35	1.40
3	B	403	NAD	O2D-C2D	2.30	1.48	1.43
3	B	403	NAD	O2B-C2B	2.33	1.48	1.43
3	B	403	NAD	C2A-N3A	2.87	1.37	1.32
3	B	403	NAD	C8A-N7A	3.16	1.40	1.34
3	B	403	NAD	C7N-N7N	5.20	1.43	1.33
3	B	403	NAD	O4D-C1D	12.24	1.56	1.41
3	B	403	NAD	O4B-C1B	12.97	1.57	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	NAD	N3A-C2A-N1A	-11.20	120.31	128.89
3	B	403	NAD	O3D-C3D-C4D	-5.75	93.81	111.05
3	B	403	NAD	N6A-C6A-N1A	-5.46	107.48	119.20
3	B	403	NAD	O7N-C7N-N7N	-3.80	117.24	122.59
3	B	403	NAD	C5B-C4B-C3B	-3.64	100.74	115.21
3	B	403	NAD	O2B-C2B-C3B	-2.94	102.27	111.83
3	B	403	NAD	C4B-O4B-C1B	-2.74	106.70	109.72
3	B	403	NAD	O4D-C1D-N1N	-2.58	105.30	108.13
3	B	403	NAD	C1B-N9A-C4A	-2.00	123.92	126.94
3	B	403	NAD	O2A-PA-O5B	2.02	118.63	108.46
3	B	403	NAD	O2A-PA-O3	2.34	115.72	105.09
3	B	403	NAD	O2N-PN-O3	2.35	115.75	105.09
3	B	403	NAD	C2N-C3N-C4N	2.57	121.16	118.29

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	B	403	NAD	O3D-C3D-C2D	2.80	120.93	111.83
3	B	403	NAD	O7N-C7N-C3N	3.53	123.43	119.59
3	B	403	NAD	C4D-O4D-C1D	3.64	113.72	109.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	375/375 (100%)	0.01	0 100 100	20, 35, 48, 60	0
1	B	375/375 (100%)	0.05	4 (1%) 82 84	21, 33, 51, 60	0
All	All	750/750 (100%)	0.03	4 (0%) 91 92	20, 34, 50, 60	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	304	PHE	3.1
1	B	313	ASN	2.9
1	B	307	HIS	2.1
1	B	303	ALA	2.1

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, 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	B	401	1/1	1.00	0.17	1.31	36,36,36,36	0
2	ZN	B	402	1/1	0.99	0.18	1.26	32,32,32,32	0
2	ZN	A	400	1/1	1.00	0.15	0.33	26,26,26,26	0
2	ZN	A	401	1/1	0.99	0.14	-0.22	38,38,38,38	0
3	NAD	B	403	44/44	0.98	0.15	-0.53	14,27,36,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.