



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:54 am GMT

PDB ID : 1AXG
Title : CRYSTAL STRUCTURE OF THE VAL203->ALA MUTANT OF LIVER ALCOHOL DEHYDROGENASE COMPLEXED WITH COFACTOR NAD AND INHIBITOR TRIFLUOROETHANOL SOLVED TO 2.5 ANGSTROM RESOLUTION
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Deposited on : 1997-10-15
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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

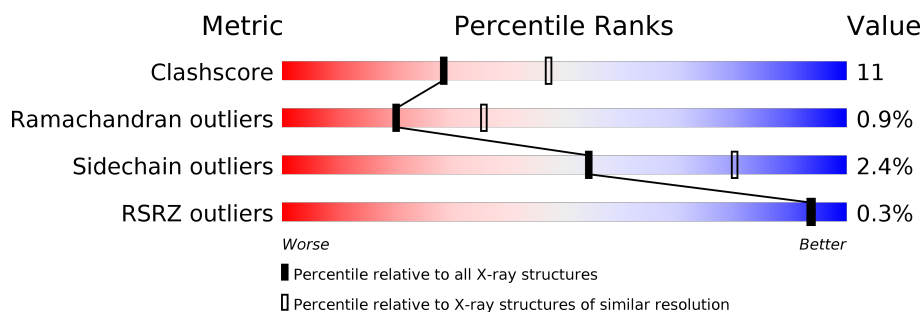
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



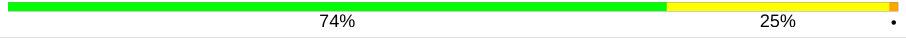

The reported resolution of this entry is 2.50 Å.

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))
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (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	374	 72% 28% .
1	B	374	 73% 26% .
1	C	374	 74% 25% .
1	D	374	 70% 29% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	A	402	-	-	X	-
4	ETF	A	404	-	-	-	X
4	ETF	B	404	-	-	-	X
4	ETF	C	404	-	-	-	X
4	ETF	D	404	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 11655 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 ALCOHOL DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	0	0	0
			2783	1767	472	521	23			
1	B	374	Total	C	N	O	S	0	0	0
			2783	1767	472	521	23			
1	C	374	Total	C	N	O	S	0	0	0
			2783	1767	472	521	23			
1	D	374	Total	C	N	O	S	0	0	0
			2783	1767	472	521	23			

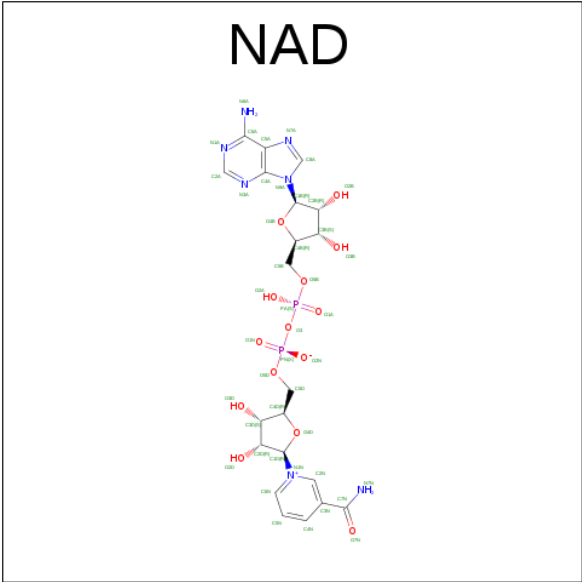
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	ALA	VAL	ENGINEERED	UNP P00327
B	203	ALA	VAL	ENGINEERED	UNP P00327
C	203	ALA	VAL	ENGINEERED	UNP P00327
D	203	ALA	VAL	ENGINEERED	UNP P00327

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

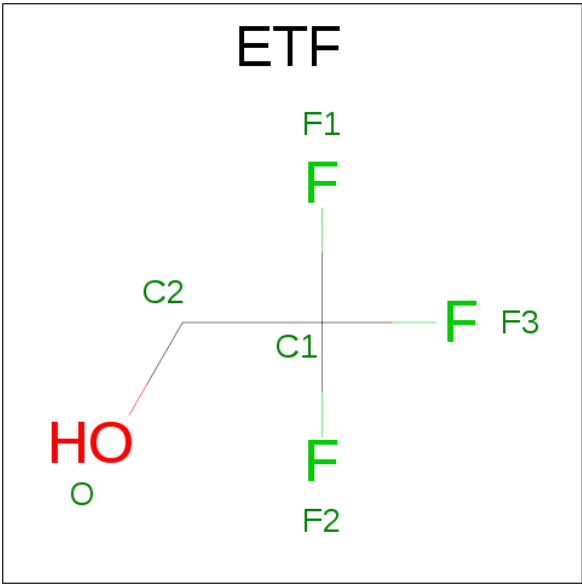
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

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



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

- Molecule 4 is TRIFLUOROETHANOL (three-letter code: ETF) (formula: C₂H₃F₃O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C F O 6 2 3 1	0	0
4	B	1	Total C F O 6 2 3 1	0	0
4	C	1	Total C F O 6 2 3 1	0	0
4	D	1	Total C F O 6 2 3 1	0	0

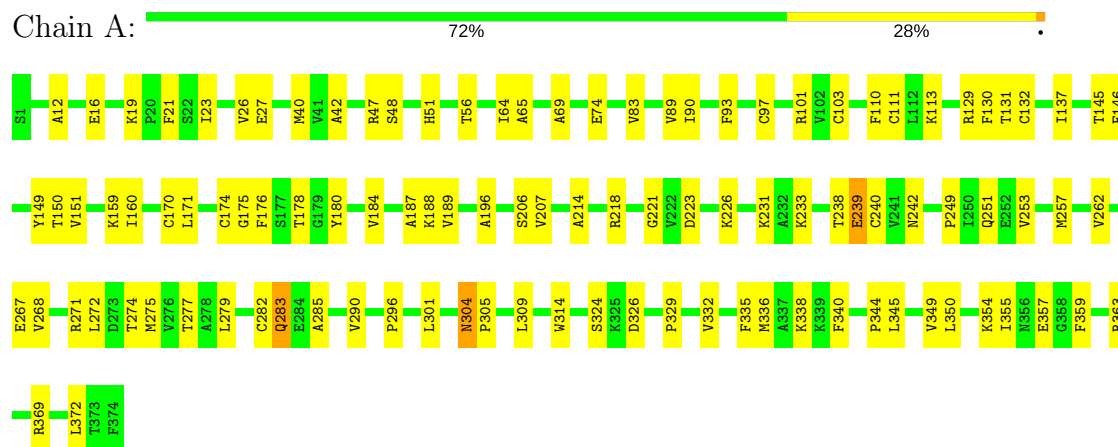
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	83	Total O 83 83	0	0
5	B	90	Total O 90 90	0	0
5	C	71	Total O 71 71	0	0
5	D	71	Total O 71 71	0	0

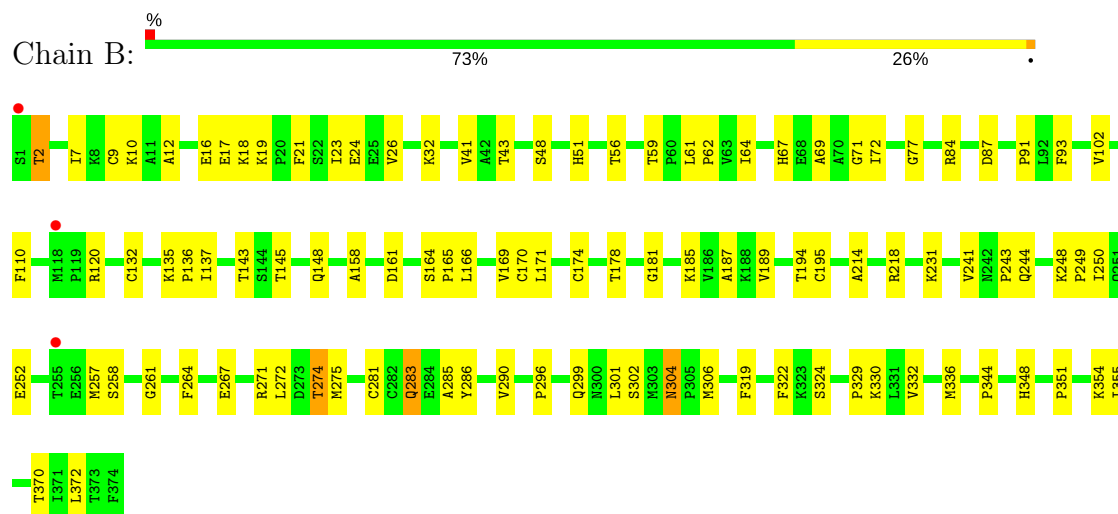
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

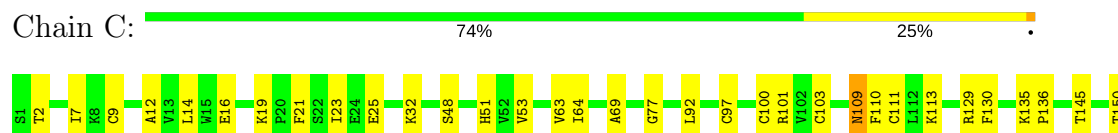
• Molecule 1: ALCOHOL DEHYDROGENASE

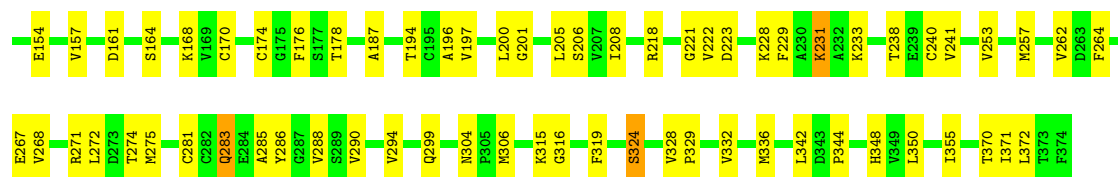


• Molecule 1: ALCOHOL DEHYDROGENASE

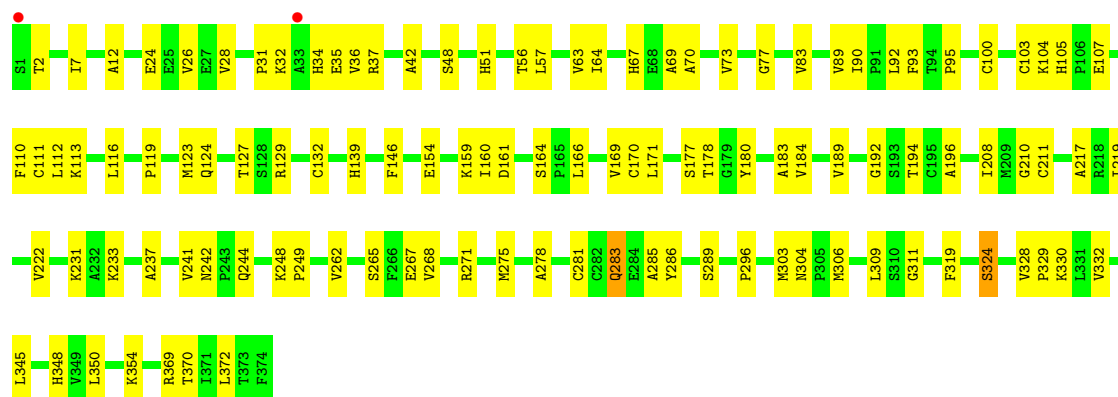


• Molecule 1: ALCOHOL DEHYDROGENASE





• Molecule 1: ALCOHOL DEHYDROGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.35Å 86.73Å 90.64Å 102.66° 99.24° 101.63°	Depositor
Resolution (Å)	10.00 – 2.50 9.99 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.0 (10.00-2.50) 78.3 (9.99-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.214 , 0.294 0.189 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.2	EDS
L-test for twinning ¹	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11655	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN, ETF, 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	0.33	0/2835	0.62	0/3831
1	B	0.33	0/2835	0.61	0/3831
1	C	0.33	0/2835	0.60	0/3831
1	D	0.33	0/2835	0.61	0/3831
All	All	0.33	0/11340	0.61	0/15324

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2783	0	2846	64	0
1	B	2783	0	2847	60	0
1	C	2783	0	2845	57	0
1	D	2783	0	2845	71	0
2	A	2	0	0	2	0
2	B	2	0	0	0	0
2	C	2	0	0	1	0
2	D	2	0	0	1	0
3	A	44	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	44	0	26	1	0
3	C	44	0	26	1	0
3	D	44	0	26	2	0
4	A	6	0	3	1	0
4	B	6	0	3	1	0
4	C	6	0	3	0	0
4	D	6	0	3	0	0
5	A	83	0	0	1	0
5	B	90	0	0	5	0
5	C	71	0	0	2	0
5	D	71	0	0	1	0
All	All	11655	0	11499	245	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 11.

All (245) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:CYS:HG	2:C:402:ZN:ZN	0.93	0.77
1:A:111:CYS:HG	2:A:402:ZN:ZN	0.99	0.72
1:C:271:ARG:HB2	1:C:274:THR:OG1	1.91	0.71
1:A:283:GLN:HE22	1:A:285:ALA:HB3	1.57	0.70
1:A:231:LYS:HE3	1:A:344:PRO:O	1.91	0.69
1:B:272:LEU:HD11	1:B:299:GLN:HB3	1.72	0.69
1:B:161:ASP:HB3	1:B:164:SER:OG	1.93	0.69
1:D:89:VAL:HG12	1:D:159:LYS:HA	1.74	0.69
1:D:2:THR:HB	1:D:7:ILE:HD11	1.74	0.69
1:A:271:ARG:HB2	1:A:274:THR:OG1	1.94	0.68
1:D:196:ALA:HB2	1:D:262:VAL:HG11	1.74	0.68
1:B:267:GLU:HG3	1:B:275:MET:HA	1.75	0.67
1:A:16:GLU:HB2	1:A:19:LYS:HG3	1.76	0.67
1:C:12:ALA:HB1	1:C:21:PHE:HB3	1.77	0.67
1:C:32:LYS:O	1:C:77:GLY:HA3	1.96	0.66
1:D:242:ASN:OD1	1:D:244:GLN:HG2	1.93	0.66
1:D:350:LEU:HD22	1:D:354:LYS:HG2	1.77	0.66
1:B:17:GLU:HG2	1:B:18:LYS:HG3	1.79	0.65
1:D:103:CYS:HG	2:D:402:ZN:ZN	1.07	0.65
1:A:180:TYR:O	1:A:184:VAL:HG22	1.97	0.64
1:D:345:LEU:O	1:D:369:ARG:HB2	1.98	0.64
1:D:92:LEU:HD13	1:D:324:SER:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:PRO:HD2	1:B:354:LYS:HD3	1.79	0.63
1:C:196:ALA:HB2	1:C:262:VAL:HG11	1.78	0.63
1:B:26:VAL:HG12	1:B:132:CYS:HB2	1.81	0.63
1:D:83:VAL:HG12	1:D:159:LYS:HB2	1.80	0.63
1:C:16:GLU:HB2	1:C:19:LYS:HG3	1.79	0.62
1:C:161:ASP:HB3	1:C:164:SER:OG	1.98	0.62
1:A:350:LEU:O	1:A:372:LEU:HA	2.00	0.62
1:C:272:LEU:HD11	1:C:299:GLN:HB3	1.80	0.61
1:B:249:PRO:HG2	1:B:252:GLU:HG3	1.81	0.61
1:C:23:ILE:HD11	1:C:355:ILE:CG2	2.31	0.61
1:C:231:LYS:HE3	1:C:344:PRO:O	2.00	0.61
1:D:31:PRO:HD3	1:D:37:ARG:HB2	1.83	0.61
1:A:89:VAL:HG12	1:A:159:LYS:HA	1.83	0.60
1:A:251:GLN:HB3	1:A:277:THR:HG23	1.83	0.60
1:A:332:VAL:O	1:A:336:MET:HG2	2.02	0.60
1:D:123:MET:SD	1:D:139:HIS:CE1	2.94	0.60
1:B:135:LYS:HD3	5:B:430:HOH:O	2.01	0.60
1:A:97:CYS:HG	1:A:111:CYS:HG	1.47	0.59
1:D:180:TYR:O	1:D:184:VAL:HG22	2.01	0.59
1:D:306:MET:SD	1:D:309:LEU:HD12	2.42	0.59
1:A:272:LEU:HD22	1:A:301:LEU:HB3	1.85	0.58
1:A:97:CYS:HG	2:A:402:ZN:ZN	1.16	0.58
1:B:93:PHE:CE2	4:B:404:ETF:H22	2.38	0.58
1:A:26:VAL:HG12	1:A:132:CYS:HB2	1.86	0.58
1:D:192:GLY:HA2	1:D:217:ALA:HB2	1.85	0.58
1:A:12:ALA:HB1	1:A:21:PHE:HB3	1.86	0.58
1:D:12:ALA:HB3	1:D:146:PHE:HB3	1.86	0.58
1:A:304:ASN:HD22	1:A:305:PRO:HD2	1.69	0.57
1:D:208:ILE:HG23	1:D:219:ILE:HG21	1.86	0.57
1:A:271:ARG:O	1:A:275:MET:HG3	2.05	0.56
1:C:9:CYS:O	1:C:25:GLU:HA	2.05	0.56
1:C:200:LEU:HG	1:C:223:ASP:HB2	1.87	0.56
1:D:194:THR:HG22	1:D:262:VAL:HG12	1.87	0.56
1:C:228:LYS:NZ	5:C:475:HOH:O	2.39	0.55
1:D:69:ALA:HA	1:D:170:CYS:HB2	1.88	0.55
1:A:97:CYS:HB3	1:A:113:LYS:HG3	1.87	0.55
1:A:83:VAL:HG12	1:A:159:LYS:HB2	1.88	0.55
1:C:290:VAL:HA	1:C:315:LYS:O	2.07	0.55
1:D:283:GLN:NE2	1:D:285:ALA:H	2.05	0.55
1:D:51:HIS:HB3	1:D:57:LEU:HB2	1.89	0.55
1:A:69:ALA:HB3	1:A:145:THR:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:ALA:HB2	1:C:290:VAL:HG21	1.89	0.55
1:A:110:PHE:HB3	1:B:285:ALA:HB2	1.90	0.54
1:B:332:VAL:O	1:B:336:MET:HG2	2.06	0.54
1:B:231:LYS:HE2	1:B:344:PRO:O	2.07	0.54
1:C:178:THR:HG21	3:C:403:NAD:C4N	2.38	0.54
1:A:188:LYS:HE3	5:A:463:HOH:O	2.06	0.54
1:B:171:LEU:HB3	5:B:478:HOH:O	2.07	0.54
1:C:97:CYS:SG	1:C:100:CYS:SG	3.06	0.54
1:B:67:HIS:HB2	1:B:93:PHE:HB3	1.89	0.53
1:A:48:SER:O	1:A:51:HIS:HB2	2.09	0.53
1:D:328:VAL:O	1:D:332:VAL:HG23	2.09	0.53
1:A:196:ALA:HB2	1:A:262:VAL:HG11	1.91	0.53
1:C:283:GLN:HE22	1:C:285:ALA:HB3	1.74	0.53
1:C:92:LEU:CD2	1:C:328:VAL:HG21	2.39	0.53
1:D:110:PHE:CE1	1:D:116:LEU:HD23	2.44	0.53
1:D:160:ILE:CG2	1:D:332:VAL:HG21	2.39	0.53
1:D:56:THR:HG23	1:D:296:PRO:HA	1.91	0.53
1:D:100:CYS:HB2	1:D:112:LEU:HD12	1.90	0.53
1:C:48:SER:O	1:C:51:HIS:HB2	2.09	0.53
1:D:271:ARG:O	1:D:275:MET:HG3	2.09	0.52
1:D:67:HIS:HB2	1:D:93:PHE:HB3	1.91	0.52
1:D:26:VAL:HG12	1:D:132:CYS:HB2	1.91	0.52
1:D:48:SER:O	1:D:51:HIS:HB2	2.08	0.52
1:B:283:GLN:NE2	1:B:285:ALA:H	2.07	0.52
1:C:253:VAL:O	1:C:257:MET:HG3	2.09	0.52
1:D:166:LEU:HA	1:D:169:VAL:HG22	1.91	0.52
1:B:69:ALA:HA	1:B:170:CYS:HB2	1.92	0.52
1:B:165:PRO:HD2	1:B:336:MET:HE3	1.92	0.52
1:A:129:ARG:HG3	1:A:151:VAL:HB	1.90	0.52
1:B:195:CYS:HA	1:B:264:PHE:O	2.09	0.52
1:B:48:SER:O	1:B:51:HIS:HB2	2.09	0.52
1:C:32:LYS:HD2	1:C:129:ARG:NH2	2.25	0.51
1:A:178:THR:HG21	3:A:403:NAD:C4N	2.41	0.51
1:A:64:ILE:HD12	1:A:130:PHE:CD2	2.45	0.51
1:A:12:ALA:HB3	1:A:146:PHE:HB3	1.91	0.51
1:A:56:THR:HG23	1:A:296:PRO:HA	1.92	0.51
1:B:10:LYS:HA	1:B:24:GLU:O	2.11	0.51
1:B:336:MET:HE2	1:B:336:MET:HA	1.92	0.51
1:B:271:ARG:O	1:B:275:MET:HG3	2.11	0.51
1:C:100:CYS:SG	1:C:103:CYS:SG	3.09	0.51
1:C:271:ARG:O	1:C:275:MET:HG3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:GLU:HG3	1:A:275:MET:HA	1.93	0.50
1:B:91:PRO:HB2	1:B:143:THR:HG22	1.93	0.50
1:C:201:GLY:O	1:C:205:LEU:HG	2.12	0.50
1:B:304:ASN:HD22	1:B:306:MET:H	1.58	0.50
1:A:338:LYS:HE2	1:D:119:PRO:O	2.10	0.50
1:D:283:GLN:HE22	1:D:285:ALA:HB3	1.76	0.50
1:A:176:PHE:HA	1:A:206:SER:OG	2.11	0.50
1:D:32:LYS:O	1:D:77:GLY:HA3	2.12	0.50
1:B:102:VAL:HG11	1:B:110:PHE:O	2.11	0.50
1:B:120:ARG:HB2	5:B:474:HOH:O	2.11	0.50
1:C:283:GLN:NE2	1:C:285:ALA:H	2.09	0.50
1:D:267:GLU:HG3	1:D:275:MET:HA	1.92	0.50
1:A:103:CYS:SG	1:A:111:CYS:SG	3.09	0.50
1:A:349:VAL:O	1:A:350:LEU:HD23	2.12	0.49
1:A:359:PHE:HB3	1:A:363:ARG:NH1	2.27	0.49
1:C:48:SER:HA	1:C:51:HIS:CD2	2.47	0.49
1:D:328:VAL:HB	1:D:329:PRO:HD3	1.94	0.49
1:B:2:THR:HB	1:B:7:ILE:HD11	1.94	0.49
1:A:90:ILE:HG13	1:A:160:ILE:HD13	1.94	0.49
1:C:176:PHE:HA	1:C:206:SER:OG	2.12	0.49
1:C:350:LEU:O	1:C:372:LEU:HA	2.13	0.49
1:D:113:LYS:HB3	1:D:124:GLN:NE2	2.28	0.49
1:C:168:LYS:O	1:C:342:LEU:HD13	2.12	0.49
1:D:69:ALA:O	1:D:90:ILE:HG23	2.13	0.48
1:B:271:ARG:HB2	1:B:274:THR:OG1	2.14	0.48
1:D:113:LYS:HB3	1:D:124:GLN:HE22	1.79	0.48
1:B:178:THR:HG21	3:B:403:NAD:C4N	2.43	0.48
1:A:23:ILE:HD11	1:A:355:ILE:HG21	1.96	0.48
1:A:187:ALA:HB2	1:A:290:VAL:HG21	1.96	0.48
1:D:177:SER:HB2	1:D:319:PHE:CE1	2.48	0.48
1:B:166:LEU:HA	1:B:169:VAL:HG22	1.95	0.47
1:A:64:ILE:HG13	1:A:137:ILE:HG21	1.97	0.47
1:D:171:LEU:HD22	1:D:369:ARG:HG3	1.96	0.47
1:B:120:ARG:HG2	5:B:453:HOH:O	2.14	0.47
1:D:222:VAL:HG22	1:D:241:VAL:HG22	1.97	0.47
1:A:326:ASP:O	1:A:329:PRO:HD2	2.15	0.47
1:A:233:LYS:HE2	1:A:239:GLU:HA	1.97	0.47
1:B:248:LYS:HB2	1:B:249:PRO:HD2	1.96	0.47
1:B:189:VAL:O	1:B:214:ALA:HB1	2.15	0.47
1:D:70:ALA:HB1	1:D:166:LEU:HD22	1.96	0.47
1:A:101:ARG:NH2	1:B:261:GLY:O	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:332:VAL:O	1:C:336:MET:HG2	2.15	0.46
1:B:43:THR:HG22	1:B:372:LEU:HB2	1.97	0.46
1:B:23:ILE:HD11	1:B:355:ILE:CG2	2.46	0.46
1:D:105:HIS:CE1	1:D:107:GLU:HB2	2.51	0.46
1:D:105:HIS:HE1	1:D:107:GLU:HB2	1.79	0.46
1:B:16:GLU:HB2	1:B:19:LYS:HG3	1.96	0.46
1:C:170:CYS:SG	1:C:371:ILE:HD12	2.56	0.45
1:C:23:ILE:HD11	1:C:355:ILE:HG22	1.96	0.45
1:A:249:PRO:HB2	1:A:251:GLN:NE2	2.30	0.45
1:B:187:ALA:HB2	1:B:290:VAL:HG21	1.97	0.45
1:C:316:GLY:O	1:D:311:GLY:HA2	2.16	0.45
1:C:154:GLU:O	1:C:157:VAL:HG12	2.17	0.45
1:D:161:ASP:HB3	1:D:164:SER:HG	1.82	0.44
1:D:268:VAL:O	3:D:403:NAD:H51N	2.17	0.44
1:A:218:ARG:HA	1:A:238:THR:HG21	1.99	0.44
1:A:354:LYS:HD2	1:A:357:GLU:OE2	2.17	0.44
1:B:69:ALA:HB3	1:B:145:THR:HG21	1.99	0.44
1:D:63:VAL:HG22	1:D:64:ILE:N	2.32	0.44
1:C:348:HIS:O	1:C:370:THR:HA	2.17	0.44
1:D:90:ILE:HG13	1:D:160:ILE:HD13	1.98	0.44
1:D:34:HIS:O	1:D:154:GLU:HB2	2.17	0.44
1:B:336:MET:HE2	1:B:336:MET:CA	2.46	0.44
1:B:218:ARG:HH12	1:B:258:SER:HA	1.83	0.44
1:B:56:THR:HG23	1:B:296:PRO:HA	2.00	0.44
1:B:84:ARG:O	1:B:87:ASP:HB2	2.18	0.44
1:C:14:LEU:HD21	1:C:53:VAL:HA	1.99	0.44
1:D:95:PRO:O	1:D:111:CYS:HB2	2.18	0.44
1:A:279:LEU:HD13	1:A:314:TRP:CG	2.52	0.44
1:D:265:SER:O	1:D:289:SER:HA	2.18	0.44
1:A:171:LEU:HD22	1:A:369:ARG:HG3	1.98	0.43
1:C:109:ASN:O	1:C:111:CYS:N	2.51	0.43
1:C:69:ALA:HB3	1:C:145:THR:HG21	1.99	0.43
1:C:222:VAL:HG22	1:C:241:VAL:CG2	2.48	0.43
1:C:264:PHE:CD1	1:C:288:VAL:HB	2.53	0.43
1:D:110:PHE:CZ	1:D:116:LEU:HD23	2.53	0.43
1:D:36:VAL:HG13	1:D:73:VAL:HG13	1.99	0.43
1:A:93:PHE:CE1	4:A:404:ETF:H22	2.53	0.43
1:B:194:THR:HA	1:B:218:ARG:O	2.18	0.43
1:B:244:GLN:HB2	5:B:434:HOH:O	2.19	0.43
1:C:113:LYS:HE3	5:C:446:HOH:O	2.17	0.43
1:D:278:ALA:HB3	5:D:473:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:100:CYS:O	1:D:104:LYS:HG3	2.19	0.43
1:C:221:GLY:O	1:C:240:CYS:HA	2.19	0.43
1:A:253:VAL:O	1:A:257:MET:HG3	2.18	0.43
1:A:74:GLU:HG3	1:A:149:TYR:HE2	1.82	0.43
1:A:262:VAL:HG22	1:A:282:CYS:HA	2.00	0.43
1:B:158:ALA:HB1	1:B:329:PRO:HG3	2.01	0.43
1:C:264:PHE:HD1	1:C:288:VAL:HB	1.83	0.43
1:D:48:SER:HA	1:D:51:HIS:CD2	2.54	0.43
1:C:294:VAL:HA	1:D:309:LEU:HD11	2.00	0.43
1:B:41:VAL:HG23	1:B:71:GLY:HA2	2.00	0.43
1:B:135:LYS:HA	1:B:136:PRO:HD3	1.86	0.43
1:B:32:LYS:O	1:B:77:GLY:HA3	2.19	0.43
1:B:348:HIS:O	1:B:370:THR:HA	2.18	0.43
1:D:42:ALA:HA	1:D:372:LEU:O	2.18	0.43
1:D:24:GLU:HB3	1:D:132:CYS:SG	2.59	0.43
1:A:47:ARG:HG3	1:A:51:HIS:CE1	2.54	0.42
1:B:72:ILE:HA	1:B:87:ASP:O	2.19	0.42
1:C:328:VAL:HB	1:C:329:PRO:HD3	2.00	0.42
1:C:63:VAL:HG22	1:C:64:ILE:N	2.34	0.42
1:D:178:THR:HG21	3:D:403:NAD:C4N	2.49	0.42
1:D:183:ALA:HB3	1:D:210:GLY:HA3	2.01	0.42
1:A:221:GLY:O	1:A:240:CYS:HA	2.18	0.42
1:B:181:GLY:O	1:B:185:LYS:HB3	2.19	0.42
1:C:2:THR:HB	1:C:7:ILE:HD11	2.00	0.42
1:A:335:PHE:HB2	1:A:340:PHE:CZ	2.55	0.42
1:A:69:ALA:HA	1:A:170:CYS:HB2	2.02	0.42
1:A:189:VAL:HB	1:A:214:ALA:CB	2.48	0.42
1:B:241:VAL:HG11	1:B:257:MET:SD	2.60	0.42
1:C:101:ARG:HE	1:D:283:GLN:HG3	1.84	0.42
1:B:62:PRO:HB2	1:B:137:ILE:HD12	2.01	0.42
1:D:283:GLN:HE22	1:D:285:ALA:H	1.66	0.42
1:D:348:HIS:O	1:D:370:THR:HA	2.19	0.42
1:C:130:PHE:HZ	1:C:150:THR:OG1	2.03	0.42
1:D:248:LYS:HB2	1:D:249:PRO:HD2	2.02	0.42
1:A:305:PRO:HB3	1:B:301:LEU:HD23	2.01	0.41
1:C:23:ILE:HD11	1:C:355:ILE:HG21	2.00	0.41
1:A:223:ASP:O	1:A:242:ASN:HA	2.20	0.41
1:A:207:VAL:HG21	1:A:268:VAL:HG21	2.02	0.41
1:A:27:GLU:HB2	1:A:131:THR:HB	2.02	0.41
1:A:345:LEU:O	1:A:369:ARG:HB2	2.21	0.41
1:B:243:PRO:HA	1:B:250:ILE:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:ALA:HB1	1:B:21:PHE:HB3	2.03	0.41
1:C:197:VAL:HG21	1:C:208:ILE:HG12	2.01	0.41
1:C:283:GLN:HE22	1:C:285:ALA:H	1.68	0.41
1:A:42:ALA:HA	1:A:372:LEU:O	2.20	0.41
1:C:194:THR:HG23	1:C:218:ARG:HB3	2.02	0.41
1:C:319:PHE:HE2	1:C:324:SER:HA	1.86	0.41
1:C:229:PHE:O	1:C:233:LYS:HG2	2.21	0.41
1:D:233:LYS:HA	1:D:237:ALA:HB3	2.02	0.41
1:B:319:PHE:CD2	1:B:322:PHE:HB2	2.55	0.41
1:C:135:LYS:HA	1:C:136:PRO:HD3	1.92	0.41
1:D:28:VAL:HA	1:D:129:ARG:O	2.20	0.41
1:D:189:VAL:HG21	1:D:211:CYS:SG	2.61	0.41
1:A:40:MET:CE	1:A:150:THR:HG22	2.51	0.41
1:B:9:CYS:HB2	1:B:148:GLN:HB2	2.03	0.41
1:A:304:ASN:HA	1:A:305:PRO:HD2	1.92	0.40
1:B:64:ILE:HG13	1:B:137:ILE:HG21	2.03	0.40
1:D:100:CYS:SG	1:D:103:CYS:SG	3.04	0.40
1:D:31:PRO:HB2	1:D:35:GLU:O	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	372/374 (100%)	339 (91%)	29 (8%)	4 (1%)	17	29
1	B	372/374 (100%)	339 (91%)	30 (8%)	3 (1%)	22	39
1	C	372/374 (100%)	341 (92%)	27 (7%)	4 (1%)	17	29
1	D	372/374 (100%)	336 (90%)	34 (9%)	2 (0%)	32	53
All	All	1488/1496 (100%)	1355 (91%)	120 (8%)	13 (1%)	20	36

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	175	GLY
1	B	174	CYS
1	C	324	SER
1	D	286	TYR
1	A	174	CYS
1	A	324	SER
1	B	286	TYR
1	C	174	CYS
1	D	324	SER
1	B	324	SER
1	C	110	PHE
1	C	286	TYR
1	A	65	ALA

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	307/307 (100%)	302 (98%)	5 (2%)	68	88
1	B	307/307 (100%)	298 (97%)	9 (3%)	48	75
1	C	307/307 (100%)	298 (97%)	9 (3%)	48	75
1	D	307/307 (100%)	300 (98%)	7 (2%)	56	81
All	All	1228/1228 (100%)	1198 (98%)	30 (2%)	54	80

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	226	LYS
1	A	239	GLU
1	A	283	GLN
1	A	304	ASN
1	A	309	LEU
1	B	2	THR
1	B	59	THR
1	B	61	LEU
1	B	274	THR

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Mol	Chain	Res	Type
1	B	281	CYS
1	B	283	GLN
1	B	302	SER
1	B	304	ASN
1	B	330	LYS
1	C	109	ASN
1	C	231	LYS
1	C	238	THR
1	C	267	GLU
1	C	268	VAL
1	C	281	CYS
1	C	283	GLN
1	C	304	ASN
1	C	306	MET
1	D	127	THR
1	D	231	LYS
1	D	281	CYS
1	D	283	GLN
1	D	303	MET
1	D	304	ASN
1	D	330	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	HIS
1	A	283	GLN
1	A	304	ASN
1	B	34	HIS
1	B	138	HIS
1	B	283	GLN
1	B	304	ASN
1	C	283	GLN
1	C	299	GLN
1	C	300	ASN
1	C	304	ASN
1	D	138	HIS
1	D	283	GLN
1	D	300	ASN
1	D	304	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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	A	403	-	41,48,48	0.98	2 (4%)	43,73,73	1.70	4 (9%)
4	ETF	A	404	2	5,5,5	0.36	0	7,7,7	0.39	0
3	NAD	B	403	-	41,48,48	0.96	3 (7%)	43,73,73	1.79	4 (9%)
4	ETF	B	404	2	5,5,5	0.34	0	7,7,7	0.44	0
3	NAD	C	403	-	41,48,48	0.93	1 (2%)	43,73,73	1.85	8 (18%)
4	ETF	C	404	2	5,5,5	0.35	0	7,7,7	0.36	0
3	NAD	D	403	-	41,48,48	1.02	4 (9%)	43,73,73	1.78	6 (13%)
4	ETF	D	404	2	5,5,5	0.46	0	7,7,7	0.45	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
3	NAD	A	403	-	-	0/22/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ETF	A	404	2	-	0/3/3/3	0/0/0/0
3	NAD	B	403	-	-	0/22/62/62	0/5/5/5
4	ETF	B	404	2	-	0/3/3/3	0/0/0/0
3	NAD	C	403	-	-	0/22/62/62	0/5/5/5
4	ETF	C	404	2	-	0/3/3/3	0/0/0/0
3	NAD	D	403	-	-	0/22/62/62	0/5/5/5
4	ETF	D	404	2	-	0/3/3/3	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	403	NAD	C5A-N7A	-2.10	1.32	1.39
3	B	403	NAD	C5A-N7A	-2.09	1.32	1.39
3	D	403	NAD	C6N-N1N	2.06	1.40	1.35
3	B	403	NAD	C3N-C7N	2.18	1.53	1.50
3	C	403	NAD	C3N-C7N	2.30	1.54	1.50
3	D	403	NAD	O4D-C1D	2.40	1.44	1.41
3	D	403	NAD	O4B-C1B	2.51	1.44	1.41
3	D	403	NAD	C3N-C7N	2.83	1.54	1.50
3	A	403	NAD	O4D-C1D	3.23	1.45	1.41
3	B	403	NAD	O4D-C1D	3.25	1.45	1.41

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	403	NAD	N3A-C2A-N1A	-8.74	121.25	128.86
3	C	403	NAD	N3A-C2A-N1A	-8.43	121.51	128.86
3	B	403	NAD	N3A-C2A-N1A	-8.40	121.54	128.86
3	A	403	NAD	N3A-C2A-N1A	-7.90	121.98	128.86
3	C	403	NAD	C1B-N9A-C4A	-2.93	121.58	126.64
3	C	403	NAD	C4B-O4B-C1B	-2.51	107.10	109.77
3	D	403	NAD	C1B-N9A-C4A	-2.16	122.90	126.64
3	C	403	NAD	O5B-C5B-C4B	-2.10	101.56	109.00
3	C	403	NAD	C3N-C7N-N7N	-2.06	115.42	117.77
3	D	403	NAD	C2A-N1A-C6A	2.02	122.31	118.77
3	D	403	NAD	N6A-C6A-N1A	2.05	122.83	118.77
3	D	403	NAD	O7N-C7N-C3N	2.13	122.11	119.62
3	C	403	NAD	C2A-N1A-C6A	2.15	122.53	118.77
3	C	403	NAD	O4B-C4B-C3B	2.25	109.64	105.17
3	A	403	NAD	C2A-N1A-C6A	2.26	122.72	118.77
3	B	403	NAD	N6A-C6A-N1A	2.42	123.56	118.77
3	A	403	NAD	N6A-C6A-N1A	2.65	124.02	118.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	NAD	O4B-C4B-C5B	2.96	119.38	109.40
3	C	403	NAD	C4A-C5A-N7A	3.53	112.82	109.41
3	A	403	NAD	C4A-C5A-N7A	4.29	113.56	109.41
3	D	403	NAD	C4A-C5A-N7A	4.30	113.57	109.41
3	B	403	NAD	C4A-C5A-N7A	4.55	113.80	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	NAD	1	0
4	A	404	ETF	1	0
3	B	403	NAD	1	0
4	B	404	ETF	1	0
3	C	403	NAD	1	0
3	D	403	NAD	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	374/374 (100%)	-0.41	0 100 100	3, 11, 24, 33	0
1	B	374/374 (100%)	-0.46	3 (0%) 86 86	2, 10, 25, 41	0
1	C	374/374 (100%)	-0.47	0 100 100	2, 11, 24, 35	0
1	D	374/374 (100%)	-0.35	2 (0%) 90 91	2, 12, 31, 43	0
All	All	1496/1496 (100%)	-0.42	5 (0%) 93 94	2, 11, 27, 43	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	SER	2.8
1	D	1	SER	2.8
1	B	255	THR	2.6
1	D	33	ALA	2.5
1	B	118	MET	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, 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(Å ²)	Q<0.9
4	ETF	C	404	6/6	0.81	0.33	12.82	-1,11,15,17	0
4	ETF	B	404	6/6	0.78	0.27	9.84	2,12,21,24	0
4	ETF	D	404	6/6	0.84	0.27	5.65	5,9,14,15	0
4	ETF	A	404	6/6	0.88	0.21	5.29	9,11,16,17	0
3	NAD	D	403	44/44	0.96	0.11	-0.45	2,6,13,22	0
3	NAD	B	403	44/44	0.97	0.10	-0.63	2,4,11,15	0
3	NAD	A	403	44/44	0.96	0.11	-0.71	2,6,15,20	0
3	NAD	C	403	44/44	0.97	0.10	-0.75	2,9,17,19	0
2	ZN	B	402	1/1	1.00	0.02	-2.12	14,14,14,14	0
2	ZN	A	402	1/1	1.00	0.06	-2.27	11,11,11,11	0
2	ZN	C	402	1/1	0.97	0.07	-2.45	33,33,33,33	0
2	ZN	C	401	1/1	0.97	0.07	-2.68	10,10,10,10	0
2	ZN	D	401	1/1	1.00	0.04	-2.84	25,25,25,25	0
2	ZN	D	402	1/1	0.99	0.03	-3.65	27,27,27,27	0
2	ZN	B	401	1/1	0.99	0.04	-5.42	17,17,17,17	0
2	ZN	A	401	1/1	0.99	0.02	-6.90	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.