



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:16 AM GMT

PDB ID : 2Q2Q  
Title : Structure of D-3-Hydroxybutyrate Dehydrogenase from Pseudomonas putida  
Authors : Paithankar, K.S.; Feller, C.; Kuettner, E.B.; Keim, A.; Grunow, M.; Strater, N.  
Deposited on : 2007-05-29  
Resolution : 2.02 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

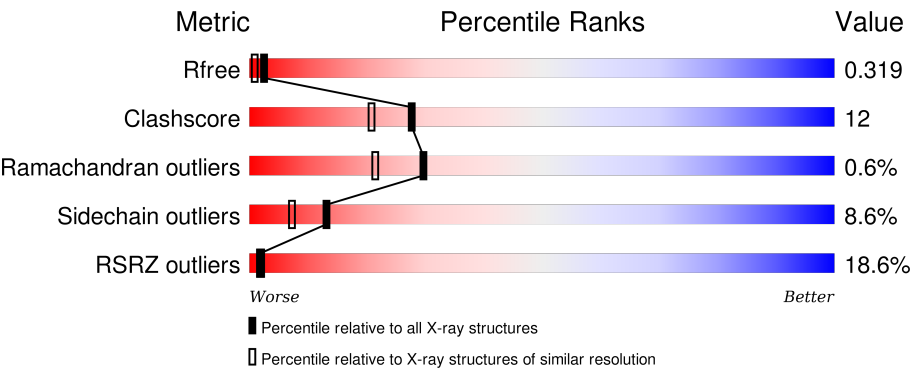
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.02 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 <sub>free</sub>	91344	7858 (2.04-2.00)
Clashscore	102246	9060 (2.04-2.00)
Ramachandran outliers	100387	8952 (2.04-2.00)
Sidechain outliers	100360	8951 (2.04-2.00)
RSRZ outliers	91569	7873 (2.04-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	255	
1	B	255	
1	C	255	
1	D	255	
1	E	255	

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Mol	Chain	Length	Quality of chain
1	F	255	<div><div>15%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>69%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>20%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>7%</div><div><div></div><div></div><div></div><div></div><div></div></div></div>
1	G	255	<div><div>19%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>73%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>22%</div><div><div></div><div></div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div></div>
1	H	255	<div><div>19%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>71%</div><div><div></div><div></div><div></div><div></div><div></div></div><div>19%</div><div><div></div><div></div><div></div><div></div><div></div></div><div><div></div><div></div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15261 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 Beta-D-hydroxybutyrate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1843	1171	330	338	4			
1	B	252	Total	C	N	O	S	0	0	0
			1852	1176	332	340	4			
1	C	243	Total	C	N	O	S	0	0	0
			1784	1137	319	324	4			
1	D	255	Total	C	N	O	S	0	0	0
			1869	1185	336	344	4			
1	E	249	Total	C	N	O	S	0	0	0
			1833	1165	329	335	4			
1	F	238	Total	C	N	O	S	0	0	0
			1742	1110	311	317	4			
1	G	251	Total	C	N	O	S	0	0	0
			1845	1173	331	337	4			
1	H	240	Total	C	N	O	S	0	0	0
			1758	1120	314	320	4			

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	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		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	61	Total	O	0	0
			61	61		
3	B	83	Total	O	0	0
			83	83		
3	C	37	Total	O	0	0
			37	37		
3	D	70	Total	O	0	0
			70	70		
3	E	55	Total	O	0	0
			55	55		
3	F	50	Total	O	0	0
			50	50		

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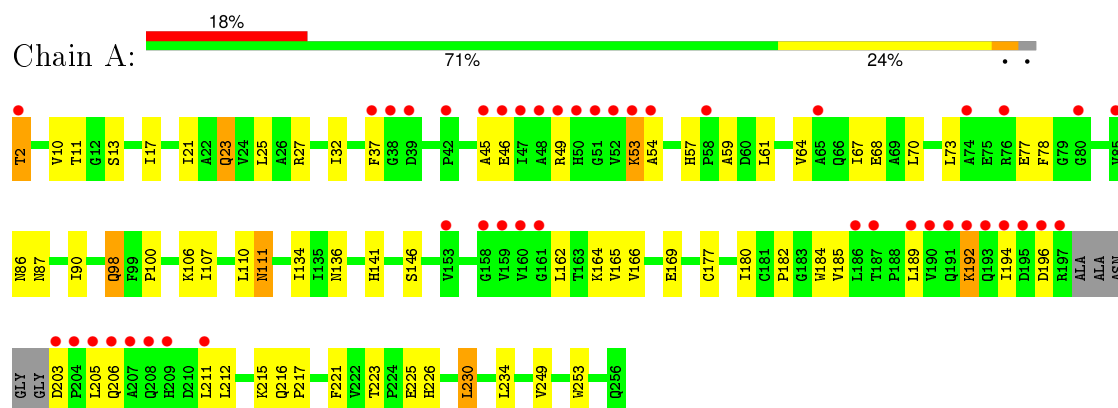
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	71	Total	O	0	0
			71	71		
3	H	44	Total	O	0	0
			44	44		

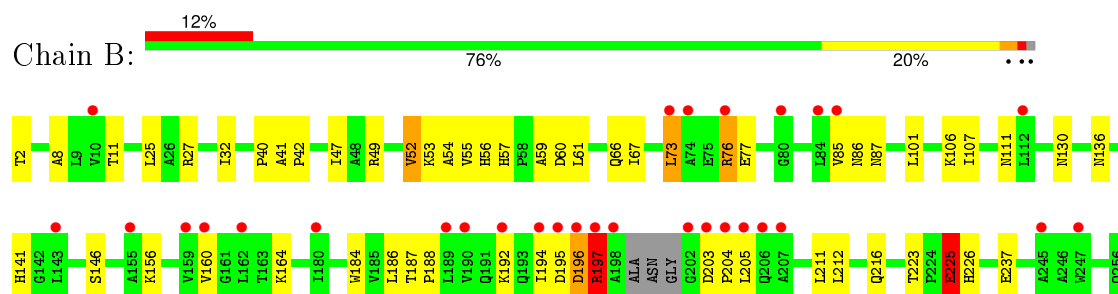
### 3 Residue-property plots

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.

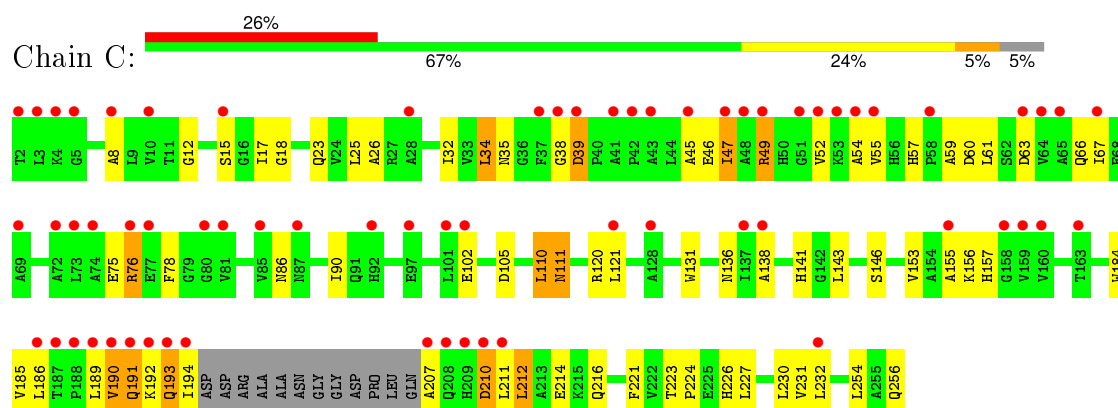
- Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



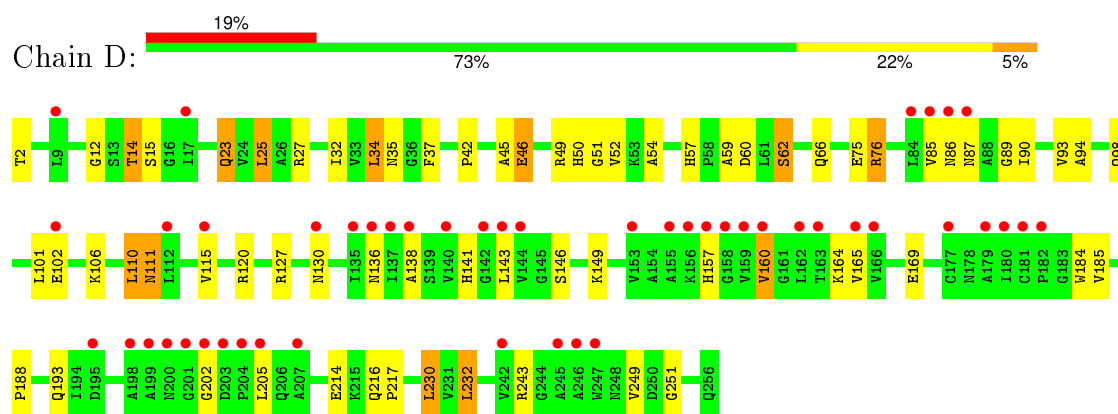
- Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



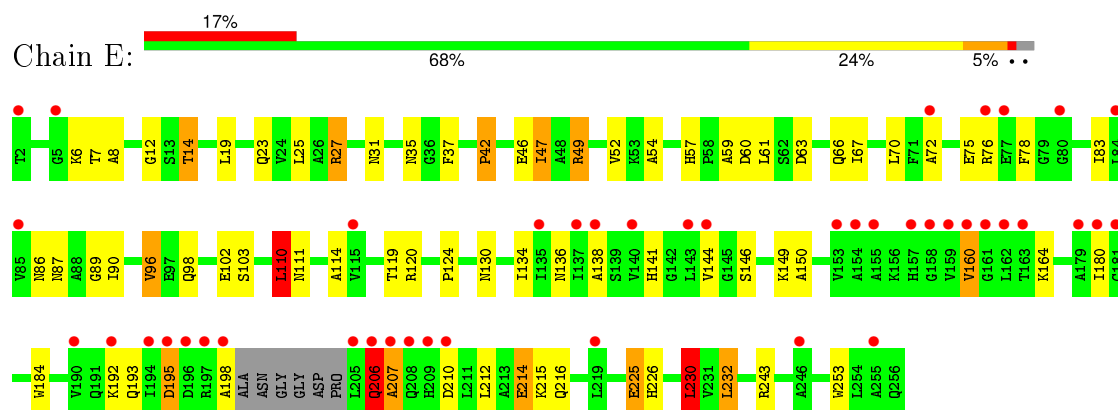
- Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



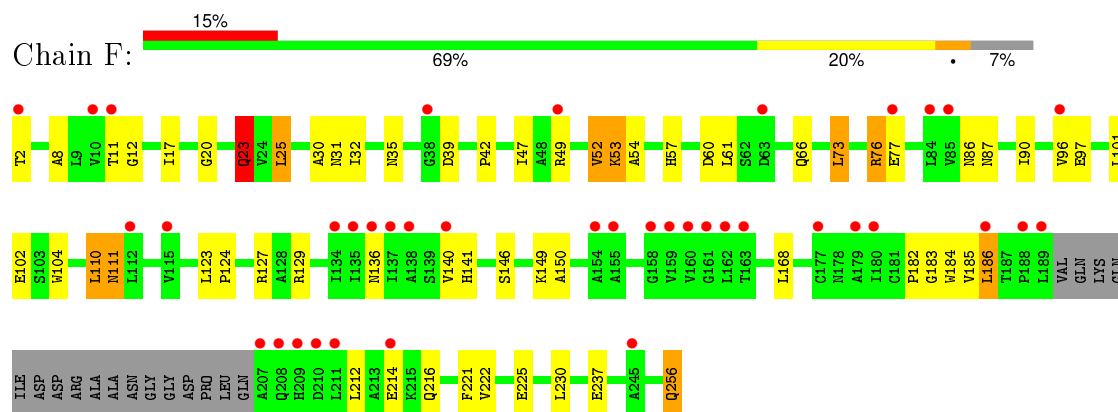
- Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



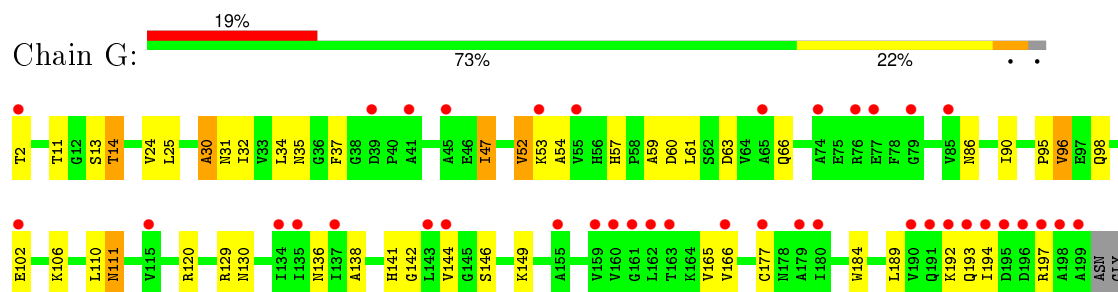
• Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



• Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



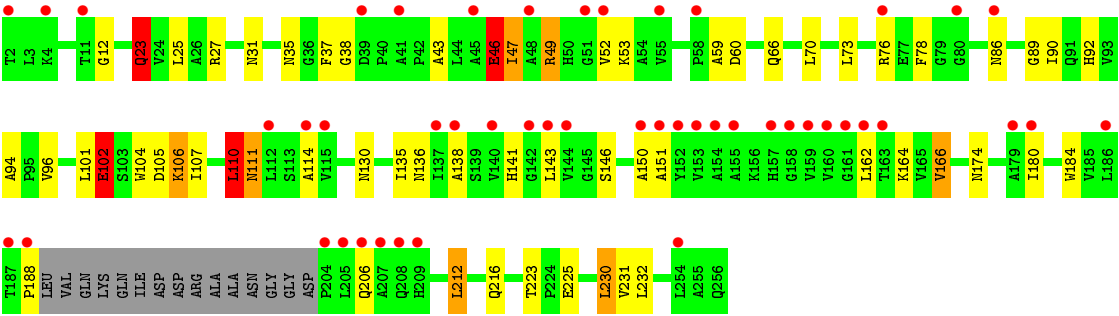
• Molecule 1: Beta-D-hydroxybutyrate dehydrogenase







● Molecule 1: Beta-D-hydroxybutyrate dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	261.46 Å   59.91 Å   116.52 Å 90.00°   113.70°   90.00°	Depositor
Resolution (Å)	30.00 – 2.02 29.96 – 2.02	Depositor EDS
% Data completeness (in resolution range)	96.4 (30.00-2.02) 96.4 (29.96-2.02)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.27 (at 2.03 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.199 , 0.274 0.255 , 0.319	Depositor DCC
$R_{free}$ test set	5241 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.6	Xtriage
Anisotropy	0.515	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 57.4	EDS
Estimated twinning fraction	0.009 for -h-2*l,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 104909 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	15261	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	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 44.55 % 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 1.5024e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.20	4/1879 (0.2%)	0.99	1/2560 (0.0%)
1	B	1.20	5/1888 (0.3%)	1.04	5/2572 (0.2%)
1	C	1.06	2/1819 (0.1%)	1.00	2/2478 (0.1%)
1	D	1.39	12/1906 (0.6%)	1.16	15/2598 (0.6%)
1	E	1.37	8/1868 (0.4%)	1.18	11/2544 (0.4%)
1	F	1.33	8/1777 (0.5%)	1.05	5/2422 (0.2%)
1	G	1.33	14/1881 (0.7%)	1.13	10/2562 (0.4%)
1	H	1.32	8/1794 (0.4%)	1.11	7/2445 (0.3%)
All	All	1.28	61/14812 (0.4%)	1.08	56/20181 (0.3%)

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	E	0	2
1	H	0	2
All	All	0	4

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	206	GLN	CD-OE1	15.60	1.58	1.24
1	G	209	HIS	CE1-NE2	14.02	1.65	1.32
1	E	206	GLN	CD-NE2	13.54	1.66	1.32
1	D	75	GLU	CG-CD	12.25	1.70	1.51
1	G	209	HIS	CG-ND1	10.21	1.61	1.38
1	D	214	GLU	CG-CD	9.54	1.66	1.51
1	F	214	GLU	CB-CG	8.57	1.68	1.52
1	F	102	GLU	CB-CG	8.56	1.68	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	23	GLN	CG-CD	8.55	1.70	1.51
1	D	115	VAL	CB-CG1	8.27	1.70	1.52
1	G	209	HIS	CG-CD2	7.61	1.48	1.35
1	F	102	GLU	CG-CD	7.44	1.63	1.51
1	F	104	TRP	CB-CG	7.31	1.63	1.50
1	D	75	GLU	CB-CG	7.25	1.66	1.52
1	D	46	GLU	CD-OE2	7.23	1.33	1.25
1	H	102	GLU	CG-CD	7.09	1.62	1.51
1	E	225	GLU	CG-CD	7.08	1.62	1.51
1	D	185	VAL	CB-CG1	6.99	1.67	1.52
1	E	160	VAL	CB-CG2	6.90	1.67	1.52
1	D	62	SER	CB-OG	-6.81	1.33	1.42
1	F	256	GLN	C-OXT	6.75	1.36	1.23
1	G	30	ALA	CA-CB	6.59	1.66	1.52
1	E	144	VAL	CB-CG2	6.59	1.66	1.52
1	G	237	GLU	CG-CD	6.52	1.61	1.51
1	B	85	VAL	CB-CG1	6.32	1.66	1.52
1	A	225	GLU	CG-CD	6.30	1.61	1.51
1	H	102	GLU	CB-CG	6.21	1.64	1.52
1	B	237	GLU	CG-CD	6.12	1.61	1.51
1	D	23	GLN	CG-CD	5.95	1.64	1.51
1	H	188	PRO	C-O	5.83	1.34	1.23
1	G	177	CYS	CB-SG	-5.74	1.72	1.81
1	H	94	ALA	CA-CB	5.67	1.64	1.52
1	F	23	GLN	CG-CD	5.64	1.64	1.51
1	E	102	GLU	CG-CD	5.63	1.60	1.51
1	C	155	ALA	CA-CB	5.61	1.64	1.52
1	G	102	GLU	CG-CD	5.61	1.60	1.51
1	A	225	GLU	CB-CG	5.60	1.62	1.52
1	E	214	GLU	CD-OE1	5.59	1.31	1.25
1	F	53	LYS	CE-NZ	5.57	1.62	1.49
1	E	96	VAL	CB-CG2	-5.54	1.41	1.52
1	B	160	VAL	CB-CG2	5.52	1.64	1.52
1	D	249	VAL	CB-CG1	5.50	1.64	1.52
1	A	98	GLN	CB-CG	-5.47	1.37	1.52
1	B	2	THR	CB-CG2	5.47	1.70	1.52
1	H	23	GLN	CB-CG	5.46	1.67	1.52
1	B	225	GLU	CG-CD	5.46	1.60	1.51
1	A	2	THR	CA-CB	5.42	1.67	1.53
1	D	76	ARG	CG-CD	5.39	1.65	1.51
1	H	166	VAL	CB-CG2	5.37	1.64	1.52
1	G	130	ASN	CB-CG	5.35	1.63	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	144	VAL	CB-CG1	5.32	1.64	1.52
1	D	160	VAL	CB-CG1	5.30	1.64	1.52
1	G	165	VAL	CA-CB	5.27	1.65	1.54
1	H	164	LYS	CD-CE	5.24	1.64	1.51
1	F	149	LYS	CD-CE	5.23	1.64	1.51
1	G	142	GLY	N-CA	5.22	1.53	1.46
1	G	166	VAL	CB-CG2	5.22	1.63	1.52
1	C	46	GLU	C-O	5.20	1.33	1.23
1	D	85	VAL	CB-CG2	5.10	1.63	1.52
1	G	96	VAL	CB-CG1	-5.06	1.42	1.52
1	G	24	VAL	CB-CG2	5.01	1.63	1.52

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	243	ARG	NE-CZ-NH2	-13.92	113.34	120.30
1	E	243	ARG	NE-CZ-NH1	12.62	126.61	120.30
1	G	129	ARG	NE-CZ-NH2	-9.81	115.40	120.30
1	C	63	ASP	CB-CG-OD2	8.27	125.74	118.30
1	E	63	ASP	CB-CG-OD1	7.68	125.21	118.30
1	D	120	ARG	NE-CZ-NH1	7.61	124.10	120.30
1	D	230	LEU	CB-CG-CD2	7.30	123.41	111.00
1	D	62	SER	CB-CA-C	-7.24	96.34	110.10
1	A	98	GLN	CB-CA-C	-7.20	96.01	110.40
1	D	243	ARG	NE-CZ-NH2	7.17	123.89	120.30
1	E	230	LEU	CB-CG-CD2	7.13	123.12	111.00
1	E	195	ASP	CB-CG-OD2	-7.02	111.98	118.30
1	H	106	LYS	CD-CE-NZ	7.02	127.85	111.70
1	D	143	LEU	CA-CB-CG	-6.94	99.34	115.30
1	H	212	LEU	CA-CB-CG	6.88	131.13	115.30
1	D	120	ARG	NE-CZ-NH2	-6.84	116.88	120.30
1	F	256	GLN	CA-C-O	-6.75	105.93	120.10
1	H	143	LEU	CB-CG-CD1	6.71	122.40	111.00
1	D	127	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	B	49	ARG	NE-CZ-NH1	6.67	123.63	120.30
1	D	34	LEU	CA-CB-CG	6.66	130.61	115.30
1	D	76	ARG	NE-CZ-NH1	6.63	123.62	120.30
1	F	127	ARG	NE-CZ-NH1	6.63	123.61	120.30
1	G	96	VAL	CB-CA-C	-6.61	98.85	111.40
1	E	96	VAL	CG1-CB-CG2	-6.47	100.55	110.90
1	F	183	GLY	N-CA-C	-6.44	96.99	113.10
1	G	205	LEU	CA-CB-CG	6.27	129.73	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	120	ARG	NE-CZ-NH2	-6.26	117.17	120.30
1	E	243	ARG	CG-CD-NE	-6.07	99.05	111.80
1	B	49	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	H	110	LEU	CB-CG-CD2	5.92	121.06	111.00
1	D	34	LEU	CB-CG-CD1	5.91	121.05	111.00
1	E	110	LEU	CB-CG-CD2	5.87	120.97	111.00
1	D	243	ARG	NE-CZ-NH1	-5.85	117.37	120.30
1	E	195	ASP	CB-CG-OD1	5.81	123.53	118.30
1	G	129	ARG	NE-CZ-NH1	5.75	123.18	120.30
1	D	25	LEU	CB-CG-CD1	5.72	120.72	111.00
1	E	232	LEU	CA-CB-CG	5.68	128.37	115.30
1	H	105	ASP	CB-CG-OD2	5.65	123.39	118.30
1	G	120	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	D	75	GLU	CA-CB-CG	5.50	125.49	113.40
1	B	76	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	G	243	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	E	243	ARG	CD-NE-CZ	5.41	131.18	123.60
1	D	232	LEU	CB-CG-CD2	5.38	120.14	111.00
1	B	27	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	F	129	ARG	NE-CZ-NH1	-5.36	117.62	120.30
1	C	105	ASP	CB-CG-OD1	5.30	123.07	118.30
1	D	52	VAL	CG1-CB-CG2	5.24	119.29	110.90
1	F	168	LEU	CB-CG-CD2	-5.23	102.11	111.00
1	H	232	LEU	CB-CG-CD2	-5.21	102.13	111.00
1	G	63	ASP	CB-CG-OD1	5.21	122.99	118.30
1	B	52	VAL	CG1-CB-CG2	5.20	119.23	110.90
1	G	189	LEU	CB-CG-CD2	-5.19	102.17	111.00
1	H	70	LEU	CA-CB-CG	5.17	127.19	115.30
1	G	210	ASP	CB-CG-OD1	5.01	122.81	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	206	GLN	Sidechain
1	E	207	ALA	Peptide
1	H	38	GLY	Peptide
1	H	49	ARG	Peptide

## 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	1843	0	1865	57	0
1	B	1852	0	1873	35	0
1	C	1784	0	1814	56	0
1	D	1869	0	1888	43	0
1	E	1833	0	1859	51	0
1	F	1742	0	1765	46	0
1	G	1845	0	1872	33	0
1	H	1758	0	1781	38	0
2	C	44	0	26	1	0
2	D	44	0	26	2	0
2	E	44	0	26	4	0
2	F	44	0	26	3	0
2	G	44	0	26	2	0
2	H	44	0	26	1	0
3	A	61	0	0	2	0
3	B	83	0	0	2	0
3	C	37	0	0	1	0
3	D	70	0	0	4	0
3	E	55	0	0	1	1
3	F	50	0	0	1	0
3	G	71	0	0	1	1
3	H	44	0	0	2	0
All	All	15261	0	14873	350	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 12.

All (350) 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:32:ILE:HG22	1:C:34:LEU:HD13	1.27	1.08
1:F:186:LEU:HD23	1:F:186:LEU:H	1.19	1.07
1:C:32:ILE:HG22	1:C:34:LEU:CD1	1.90	1.01
1:E:184:TRP:H	1:E:216:GLN:HE22	1.09	1.00
1:C:32:ILE:CG2	1:C:34:LEU:CD1	2.39	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:184:TRP:H	1:F:216:GLN:HE22	1.08	0.98
1:F:141:HIS:HE1	1:F:146:SER:OG	1.49	0.94
1:H:184:TRP:H	1:H:216:GLN:HE22	1.10	0.94
1:D:14:THR:HG21	1:D:37:PHE:O	1.68	0.92
1:F:186:LEU:H	1:F:186:LEU:CD2	1.83	0.90
1:C:32:ILE:CG2	1:C:34:LEU:HD13	2.03	0.87
1:D:42:PRO:O	1:D:46:GLU:HG2	1.74	0.87
1:F:256:GLN:OXT	3:F:317:HOH:O	1.93	0.87
1:H:96:VAL:HG12	1:H:150:ALA:HB2	1.57	0.87
1:A:184:TRP:H	1:A:216:GLN:HE22	1.17	0.87
1:C:184:TRP:H	1:C:216:GLN:HE22	1.23	0.86
1:G:184:TRP:H	1:G:216:GLN:HE22	1.24	0.86
1:E:27:ARG:NH1	3:E:342:HOH:O	2.09	0.86
1:B:184:TRP:H	1:B:216:GLN:HE22	1.25	0.84
1:G:141:HIS:HE1	1:G:146:SER:OG	1.62	0.83
1:A:189:LEU:O	1:A:189:LEU:HD12	1.79	0.83
1:G:14:THR:HG21	1:G:37:PHE:O	1.77	0.82
1:C:45:ALA:O	1:C:49:ARG:HG2	1.80	0.81
1:A:57:HIS:CE1	1:A:59:ALA:HB2	2.15	0.81
1:D:184:TRP:H	1:D:216:GLN:HE22	1.27	0.80
1:C:32:ILE:HG21	1:C:34:LEU:CD1	2.12	0.80
1:D:51:GLY:HA2	1:F:53:LYS:HD3	1.65	0.79
1:G:57:HIS:HD2	1:G:59:ALA:H	1.31	0.79
1:E:86:ASN:HD22	1:E:136:ASN:HD21	1.32	0.78
3:D:365:HOH:O	1:H:92:HIS:HD2	1.65	0.78
1:E:90:ILE:H	1:E:111:ASN:HD21	1.33	0.77
1:C:32:ILE:HG21	1:C:34:LEU:HD11	1.68	0.76
1:G:95:PRO:HG2	1:G:98:GLN:OE1	1.86	0.75
1:A:53:LYS:HG3	1:A:78:PHE:HE1	1.52	0.75
1:E:14:THR:HG22	2:E:300:NAD:O3B	1.87	0.74
1:H:90:ILE:H	1:H:111:ASN:HD21	1.33	0.74
1:A:111:ASN:HD22	1:A:111:ASN:N	1.83	0.74
1:E:184:TRP:N	1:E:216:GLN:HE22	1.85	0.74
1:F:141:HIS:CE1	1:F:146:SER:OG	2.38	0.74
1:B:86:ASN:HD22	1:B:136:ASN:HD21	1.34	0.74
1:C:76:ARG:HA	1:C:76:ARG:NE	2.03	0.73
1:B:86:ASN:HD22	1:B:136:ASN:ND2	1.85	0.73
1:F:90:ILE:H	1:F:111:ASN:HD21	1.34	0.73
1:E:72:ALA:O	1:E:76:ARG:HG3	1.88	0.73
1:G:57:HIS:CD2	1:G:59:ALA:H	2.06	0.72
1:G:47:ILE:HD11	1:G:54:ALA:HB2	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:HIS:HB2	1:B:73:LEU:HD12	1.70	0.71
1:B:101:LEU:HD12	1:B:101:LEU:H	1.54	0.71
1:F:186:LEU:HD23	1:F:186:LEU:N	1.99	0.71
1:C:38:GLY:O	1:C:39:ASP:CB	2.38	0.71
1:G:86:ASN:HD22	1:G:136:ASN:ND2	1.88	0.71
1:D:90:ILE:H	1:D:111:ASN:HD21	1.38	0.70
1:A:184:TRP:H	1:A:216:GLN:NE2	1.90	0.69
1:E:14:THR:HG21	1:E:37:PHE:O	1.91	0.69
1:A:86:ASN:HD22	1:A:136:ASN:ND2	1.91	0.69
1:E:96:VAL:HG22	1:E:150:ALA:HB2	1.74	0.69
1:A:57:HIS:HB2	1:A:73:LEU:HD23	1.73	0.69
1:A:141:HIS:HE1	1:A:146:SER:OG	1.76	0.69
1:G:90:ILE:H	1:G:111:ASN:HD21	1.42	0.68
1:F:73:LEU:HD22	1:F:77:GLU:HG2	1.75	0.68
1:B:164:LYS:NZ	1:C:157:HIS:HE1	1.92	0.68
1:C:111:ASN:HD22	1:C:111:ASN:N	1.91	0.68
1:F:30:ALA:O	1:F:52:VAL:HG11	1.94	0.67
1:F:184:TRP:N	1:F:216:GLN:HE22	1.87	0.67
1:H:90:ILE:HG13	1:H:107:ILE:CD1	2.25	0.67
1:E:86:ASN:HD22	1:E:136:ASN:ND2	1.93	0.67
1:H:184:TRP:N	1:H:216:GLN:HE22	1.90	0.66
1:D:12:GLY:H	1:D:35:ASN:HD22	1.42	0.66
1:G:86:ASN:HD22	1:G:136:ASN:HD21	1.42	0.66
1:C:60:ASP:H	1:C:66:GLN:NE2	1.93	0.65
1:F:47:ILE:CG2	1:F:54:ALA:HB2	2.25	0.65
1:F:186:LEU:HD22	1:F:221:PHE:HB3	1.78	0.65
1:C:57:HIS:HD2	1:C:59:ALA:H	1.45	0.65
1:H:60:ASP:H	1:H:66:GLN:NE2	1.95	0.65
3:D:365:HOH:O	1:H:92:HIS:CD2	2.46	0.64
1:B:55:VAL:HG11	1:B:77:GLU:HG3	1.79	0.64
1:E:49:ARG:C	1:E:49:ARG:HD3	2.17	0.64
1:F:184:TRP:H	1:F:216:GLN:NE2	1.89	0.63
1:D:138:ALA:O	2:D:300:NAD:H6N	1.98	0.63
1:E:47:ILE:CG1	1:E:54:ALA:HB2	2.28	0.63
1:H:141:HIS:HE1	1:H:146:SER:OG	1.80	0.63
1:E:184:TRP:H	1:E:216:GLN:NE2	1.88	0.63
1:A:90:ILE:HG22	1:A:111:ASN:HD21	1.63	0.62
1:C:184:TRP:N	1:C:216:GLN:HE22	1.93	0.62
1:E:47:ILE:HD11	1:E:54:ALA:HB2	1.79	0.62
1:H:73:LEU:HD12	1:H:76:ARG:NH2	2.15	0.62
1:D:45:ALA:O	1:D:49:ARG:HD2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:143:LEU:HD12	1:C:254:LEU:CD1	2.29	0.62
1:G:256:GLN:OXT	3:G:310:HOH:O	2.16	0.62
1:A:90:ILE:HB	1:A:110:LEU:HD23	1.82	0.61
1:D:51:GLY:HA2	1:F:53:LYS:CD	2.29	0.61
1:H:102:GLU:O	1:H:106:LYS:HE3	2.01	0.61
1:A:86:ASN:HD22	1:A:136:ASN:HD21	1.47	0.61
1:C:61:LEU:HD22	1:C:67:ILE:HG12	1.83	0.61
1:F:47:ILE:HG22	1:F:54:ALA:HB2	1.81	0.60
1:H:86:ASN:HD22	1:H:136:ASN:HD21	1.50	0.60
1:D:141:HIS:HE1	1:D:146:SER:OG	1.83	0.60
1:E:138:ALA:O	2:E:300:NAD:H6N	2.02	0.60
1:D:57:HIS:CD2	1:D:59:ALA:HB2	2.36	0.60
1:A:164:LYS:NZ	1:D:157:HIS:HE1	2.00	0.60
1:D:60:ASP:OD1	1:D:62:SER:HB2	2.02	0.60
1:H:92:HIS:HE1	3:H:338:HOH:O	1.83	0.59
1:A:180:ILE:HG12	1:A:230:LEU:HD13	1.84	0.59
1:G:47:ILE:CD1	1:G:52:VAL:HG12	2.33	0.59
1:H:102:GLU:C	1:H:106:LYS:HE3	2.23	0.59
1:A:184:TRP:N	1:A:216:GLN:HE22	1.95	0.59
1:F:31:ASN:HA	1:F:52:VAL:HG13	1.85	0.59
1:C:110:LEU:HB3	1:C:111:ASN:HD22	1.68	0.59
1:H:184:TRP:H	1:H:216:GLN:NE2	1.92	0.58
1:D:98:GLN:HA	1:D:98:GLN:HE21	1.66	0.58
1:C:47:ILE:HG22	1:C:54:ALA:HB2	1.84	0.58
1:B:61:LEU:O	1:B:67:ILE:HD11	2.04	0.58
1:C:38:GLY:O	1:C:39:ASP:HB3	2.03	0.58
1:D:98:GLN:HA	1:D:98:GLN:NE2	2.19	0.58
1:E:86:ASN:HB2	1:E:136:ASN:HD22	1.67	0.58
1:A:37:PHE:HB2	3:A:315:HOH:O	2.03	0.58
1:D:149:LYS:NZ	1:D:193:GLN:HE22	2.02	0.58
1:E:31:ASN:OD1	1:E:52:VAL:HG22	2.04	0.58
1:H:86:ASN:HD22	1:H:136:ASN:ND2	2.00	0.57
1:F:12:GLY:H	1:F:35:ASN:HD22	1.52	0.57
1:H:31:ASN:HA	1:H:52:VAL:HG13	1.86	0.57
1:G:206:GLN:OE1	1:G:206:GLN:HA	2.04	0.57
1:H:225:GLU:CD	1:H:225:GLU:H	2.08	0.57
1:B:184:TRP:N	1:B:216:GLN:HE22	1.97	0.57
1:B:141:HIS:HE1	1:B:146:SER:OG	1.87	0.57
1:H:12:GLY:H	1:H:35:ASN:HD22	1.51	0.57
1:D:98:GLN:CA	1:D:98:GLN:NE2	2.67	0.56
1:E:8:ALA:HA	1:E:83:ILE:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:46:GLU:O	1:H:49:ARG:HG3	2.05	0.56
1:E:149:LYS:NZ	1:E:193:GLN:HE22	2.03	0.56
1:B:60:ASP:H	1:B:66:GLN:NE2	2.03	0.56
1:D:86:ASN:HD22	1:D:136:ASN:ND2	2.03	0.56
1:F:35:ASN:HD21	2:F:300:NAD:H2A	1.71	0.56
1:C:8:ALA:HB3	1:C:32:ILE:HD13	1.88	0.55
1:E:110:LEU:HB3	1:E:111:ASN:HD22	1.70	0.55
1:H:180:ILE:HG12	1:H:230:LEU:HD13	1.88	0.55
1:E:89:GLY:HA2	1:E:110:LEU:HD13	1.88	0.55
1:C:143:LEU:HD12	1:C:254:LEU:HD13	1.89	0.55
1:A:46:GLU:OE2	1:A:49:ARG:NH1	2.40	0.55
1:C:17:ILE:HD11	1:C:185:VAL:HG21	1.88	0.55
1:G:214:GLU:HB3	1:G:215:LYS:HG2	1.89	0.55
1:C:12:GLY:H	1:C:35:ASN:HD22	1.55	0.54
1:F:17:ILE:HD11	1:F:185:VAL:HG21	1.88	0.54
1:A:45:ALA:O	1:A:49:ARG:HG2	2.08	0.54
1:D:57:HIS:HD2	1:D:59:ALA:H	1.56	0.54
1:B:225:GLU:CD	1:B:225:GLU:H	2.10	0.54
1:H:43:ALA:O	1:H:47:ILE:HG12	2.08	0.54
1:A:100:PRO:HB3	1:H:37:PHE:CD2	2.43	0.53
1:D:111:ASN:N	1:D:111:ASN:HD22	2.07	0.53
1:B:164:LYS:HZ1	1:C:157:HIS:HE1	1.55	0.53
1:A:90:ILE:HG23	1:A:107:ILE:HG12	1.90	0.53
1:E:47:ILE:CD1	1:E:54:ALA:HB2	2.39	0.53
1:B:184:TRP:H	1:B:216:GLN:NE2	2.02	0.53
1:E:207:ALA:O	1:E:210:ASP:HB2	2.09	0.53
1:A:53:LYS:HG3	1:A:78:PHE:CE1	2.38	0.53
1:H:89:GLY:HA2	1:H:110:LEU:HD13	1.91	0.52
1:C:141:HIS:HE1	1:C:146:SER:OG	1.92	0.52
1:A:23:GLN:NE2	1:A:46:GLU:HG3	2.24	0.52
1:H:223:THR:HB	1:H:225:GLU:OE1	2.09	0.52
1:C:184:TRP:H	1:C:216:GLN:NE2	2.00	0.52
1:E:42:PRO:O	1:E:46:GLU:HG2	2.09	0.52
1:A:111:ASN:N	1:A:111:ASN:ND2	2.56	0.52
1:C:138:ALA:HA	1:C:156:LYS:HD2	1.92	0.51
1:E:206:GLN:CD	1:E:206:GLN:C	2.69	0.51
1:C:256:GLN:O	3:C:332:HOH:O	2.19	0.51
1:C:184:TRP:CD1	1:C:212:LEU:HD12	2.45	0.51
1:D:86:ASN:HD22	1:D:136:ASN:HD21	1.57	0.51
1:C:26:ALA:HB2	1:C:32:ILE:HG13	1.93	0.51
1:H:135:ILE:HD12	1:H:231:VAL:HG13	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:TRP:CZ3	1:A:211:LEU:HD21	2.46	0.51
1:A:57:HIS:HB3	1:A:70:LEU:HD13	1.92	0.51
1:F:90:ILE:HG12	1:F:110:LEU:HD12	1.93	0.51
1:B:107:ILE:HG23	1:B:111:ASN:HD22	1.75	0.51
1:H:53:LYS:HG2	1:H:78:PHE:CE2	2.46	0.51
1:D:98:GLN:CA	1:D:98:GLN:HE21	2.24	0.50
1:C:17:ILE:HA	1:C:224:PRO:HB3	1.92	0.50
1:F:47:ILE:HG21	1:F:54:ALA:HB2	1.92	0.50
1:A:61:LEU:O	1:A:67:ILE:HD11	2.12	0.50
1:E:141:HIS:HE1	1:E:146:SER:OG	1.94	0.50
1:G:11:THR:HB	1:G:61:LEU:HD11	1.93	0.50
1:D:165:VAL:O	1:D:169:GLU:HG3	2.11	0.50
1:E:19:LEU:O	1:E:23:GLN:HG3	2.12	0.49
1:D:184:TRP:H	1:D:216:GLN:NE2	2.03	0.49
1:A:141:HIS:CE1	1:A:146:SER:OG	2.62	0.49
1:H:73:LEU:HD12	1:H:76:ARG:HH21	1.75	0.49
1:C:138:ALA:O	2:C:300:NAD:H6N	2.12	0.49
1:C:191:GLN:HA	1:C:194:ILE:CD1	2.42	0.49
1:D:32:ILE:O	1:D:54:ALA:HA	2.12	0.49
1:A:23:GLN:NE2	1:A:46:GLU:O	2.45	0.49
1:E:207:ALA:HA	1:E:210:ASP:OD1	2.12	0.49
1:C:55:VAL:CG2	1:C:78:PHE:HE2	2.25	0.49
1:C:141:HIS:HB3	1:C:153:VAL:HG22	1.94	0.49
1:E:57:HIS:CD2	1:E:59:ALA:HB2	2.48	0.49
1:A:64:VAL:O	1:A:68:GLU:HG3	2.13	0.49
1:C:207:ALA:O	1:C:210:ASP:HB2	2.12	0.49
1:G:184:TRP:N	1:G:216:GLN:HE22	2.02	0.49
1:G:14:THR:HG22	2:G:300:NAD:O3B	2.13	0.49
1:E:87:ASN:O	2:E:300:NAD:H4D	2.13	0.49
1:D:86:ASN:HB2	1:D:136:ASN:HD22	1.77	0.48
1:D:27:ARG:HG3	1:D:50:HIS:CE1	2.48	0.48
1:B:40:PRO:HB3	1:B:56:HIS:CG	2.48	0.48
1:E:195:ASP:O	1:E:198:ALA:HB3	2.13	0.48
1:A:194:ILE:HD11	1:A:221:PHE:CZ	2.49	0.48
1:A:73:LEU:HD12	1:A:77:GLU:CG	2.43	0.48
1:A:57:HIS:HB2	1:A:73:LEU:CD2	2.43	0.48
1:E:214:GLU:HG3	1:E:215:LYS:HG2	1.96	0.48
1:E:31:ASN:OD1	1:E:52:VAL:CG2	2.62	0.48
1:G:57:HIS:CD2	1:G:59:ALA:HB2	2.49	0.48
1:A:192:LYS:O	1:A:196:ASP:OD2	2.32	0.48
1:C:191:GLN:HA	1:C:194:ILE:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ALA:HB3	1:B:42:PRO:HD3	1.96	0.47
1:A:73:LEU:HD12	1:A:77:GLU:HG2	1.94	0.47
1:A:57:HIS:HE1	1:A:59:ALA:HB2	1.71	0.47
1:D:57:HIS:CD2	1:D:59:ALA:H	2.33	0.47
1:G:30:ALA:O	1:G:52:VAL:HG11	2.15	0.47
1:F:20:GLY:O	1:F:23:GLN:NE2	2.47	0.47
1:C:75:GLU:OE1	1:C:76:ARG:NH1	2.47	0.47
1:G:86:ASN:HB2	1:G:136:ASN:HD22	1.80	0.47
1:D:46:GLU:OE2	1:D:49:ARG:NH1	2.47	0.47
1:C:184:TRP:HD1	1:C:212:LEU:HD12	1.79	0.47
1:F:73:LEU:HD23	1:F:76:ARG:NH2	2.29	0.47
1:F:185:VAL:HG13	1:F:222:VAL:O	2.15	0.47
1:G:194:ILE:HG21	1:G:208:GLN:HB3	1.95	0.47
1:F:39:ASP:O	1:F:42:PRO:HD2	2.13	0.47
1:F:11:THR:HB	1:F:61:LEU:HD11	1.97	0.47
1:D:160:VAL:HG12	1:D:164:LYS:HE3	1.96	0.47
1:F:86:ASN:HD22	1:F:136:ASN:ND2	2.11	0.47
1:B:223:THR:H	1:B:226:HIS:CD2	2.33	0.47
1:C:111:ASN:N	1:C:111:ASN:ND2	2.62	0.47
1:C:12:GLY:O	1:C:18:GLY:HA3	2.15	0.47
1:A:86:ASN:HB2	1:A:136:ASN:HD22	1.80	0.46
1:D:87:ASN:O	2:D:300:NAD:H4D	2.16	0.46
1:E:226:HIS:HE1	1:F:237:GLU:OE1	1.98	0.46
1:A:32:ILE:O	1:A:54:ALA:HA	2.15	0.46
1:H:130:ASN:O	1:H:174:ASN:HB2	2.15	0.46
1:G:197:ARG:NH2	1:G:214:GLU:OE2	2.42	0.46
1:G:149:LYS:NZ	1:G:193:GLN:HE22	2.13	0.46
1:A:215:LYS:HB3	1:A:253:TRP:CE2	2.50	0.46
1:G:61:LEU:HG	2:G:300:NAD:H2A	1.98	0.46
1:G:47:ILE:CD1	1:G:54:ALA:HB2	2.44	0.46
1:C:190:VAL:O	1:C:193:GLN:N	2.49	0.46
1:C:227:LEU:O	1:C:231:VAL:HG23	2.14	0.46
1:E:47:ILE:CD1	1:E:52:VAL:HG13	2.46	0.46
1:A:192:LYS:HE3	1:A:192:LYS:N	2.30	0.46
1:A:17:ILE:HG22	1:A:21:ILE:HD12	1.97	0.46
1:A:90:ILE:HG22	1:A:111:ASN:ND2	2.30	0.46
1:F:87:ASN:O	2:F:300:NAD:H4D	2.16	0.46
1:A:141:HIS:HD2	3:A:316:HOH:O	1.99	0.46
1:D:57:HIS:HD2	1:D:59:ALA:HB2	1.79	0.45
1:F:60:ASP:H	1:F:66:GLN:NE2	2.14	0.45
1:E:57:HIS:HD2	1:E:59:ALA:H	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ILE:HG12	1:A:185:VAL:HG11	1.99	0.45
1:A:90:ILE:H	1:A:111:ASN:HD21	1.63	0.45
1:C:143:LEU:CD1	1:C:254:LEU:HD13	2.46	0.45
1:F:96:VAL:HG12	1:F:150:ALA:HB2	1.99	0.45
1:C:86:ASN:HD22	1:C:136:ASN:ND2	2.14	0.45
1:G:31:ASN:OD1	1:G:53:LYS:HB2	2.17	0.45
1:H:162:LEU:O	1:H:166:VAL:HG23	2.17	0.45
1:A:73:LEU:CD1	1:A:77:GLU:HG3	2.46	0.45
1:E:61:LEU:HG	2:E:300:NAD:H2A	1.99	0.45
1:B:101:LEU:HD21	1:C:120:ARG:CZ	2.47	0.45
1:B:156:LYS:NZ	3:B:334:HOH:O	2.39	0.45
1:A:134:ILE:O	1:A:177:CYS:HA	2.16	0.45
1:E:60:ASP:H	1:E:66:GLN:NE2	2.15	0.45
1:E:49:ARG:C	1:E:49:ARG:CD	2.85	0.45
1:E:180:ILE:HG12	1:E:230:LEU:HD13	1.98	0.45
1:A:203:ASP:HB3	1:A:206:GLN:HB3	1.99	0.45
1:E:206:GLN:CD	1:E:207:ALA:N	2.70	0.44
1:F:86:ASN:HB2	1:F:136:ASN:HD22	1.82	0.44
1:F:73:LEU:HD23	1:F:76:ARG:HH21	1.82	0.44
1:B:61:LEU:HD22	1:B:67:ILE:HG12	1.99	0.44
1:D:27:ARG:CG	1:D:50:HIS:CE1	3.00	0.44
1:B:211:LEU:HA	1:B:211:LEU:HD12	1.83	0.44
1:D:184:TRP:N	1:D:216:GLN:HE22	2.05	0.44
1:B:187:THR:HB	1:B:188:PRO:CD	2.47	0.44
1:E:57:HIS:HB3	1:E:70:LEU:HD13	1.99	0.44
1:E:12:GLY:H	1:E:35:ASN:HD22	1.64	0.44
1:B:164:LYS:HZ2	1:C:157:HIS:HE1	1.62	0.44
1:C:38:GLY:O	1:C:39:ASP:HB2	2.16	0.44
1:H:23:GLN:OE1	1:H:27:ARG:NH2	2.51	0.44
1:F:225:GLU:H	1:F:225:GLU:CD	2.21	0.44
1:C:75:GLU:HG3	1:C:76:ARG:HD2	2.00	0.44
1:A:234:LEU:HD23	1:A:234:LEU:HA	1.83	0.44
1:F:61:LEU:HB2	1:F:110:LEU:HD23	2.00	0.43
1:B:73:LEU:CD2	1:B:77:GLU:HG2	2.49	0.43
1:F:186:LEU:HD22	1:F:221:PHE:CB	2.45	0.43
1:H:90:ILE:HG13	1:H:107:ILE:HD12	2.00	0.43
1:E:216:GLN:HE21	1:E:253:TRP:HE3	1.65	0.43
1:B:11:THR:O	1:B:87:ASN:HB3	2.18	0.43
1:A:162:LEU:O	1:A:166:VAL:HG23	2.18	0.43
1:H:104:TRP:CE3	1:H:151:ALA:HB2	2.54	0.43
1:A:90:ILE:CB	1:A:110:LEU:HD23	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:LEU:HD23	1:B:212:LEU:N	2.34	0.43
1:G:141:HIS:CE1	1:G:146:SER:OG	2.54	0.43
1:D:93:VAL:O	1:D:94:ALA:HB2	2.19	0.43
1:B:47:ILE:HG21	1:B:54:ALA:HB2	2.02	0.42
1:D:45:ALA:O	1:D:49:ARG:CD	2.66	0.42
1:H:23:GLN:HE22	1:H:225:GLU:HG3	1.84	0.42
1:D:60:ASP:H	1:D:66:GLN:NE2	2.18	0.42
1:A:223:THR:H	1:A:226:HIS:CD2	2.36	0.42
1:H:138:ALA:O	2:H:300:NAD:H6N	2.20	0.42
1:H:59:ALA:HA	1:H:66:GLN:HE21	1.84	0.42
1:F:86:ASN:HD22	1:F:136:ASN:HD21	1.68	0.42
1:B:196:ASP:O	1:B:197:ARG:HG2	2.19	0.42
1:F:8:ALA:HB3	1:F:32:ILE:HD13	2.00	0.42
1:D:89:GLY:HA2	1:D:110:LEU:HD13	2.02	0.42
1:F:12:GLY:H	1:F:35:ASN:ND2	2.18	0.42
1:F:25:LEU:HA	1:F:25:LEU:HD12	1.93	0.41
1:A:165:VAL:O	1:A:169:GLU:HG3	2.20	0.41
1:D:130:ASN:ND2	3:D:360:HOH:O	2.53	0.41
1:A:11:THR:O	1:A:87:ASN:HB3	2.19	0.41
1:B:59:ALA:CB	1:B:66:GLN:O	2.68	0.41
1:D:217:PRO:HD2	1:D:251:GLY:O	2.21	0.41
1:G:47:ILE:O	1:G:47:ILE:HD12	2.21	0.41
1:E:160:VAL:HG12	1:E:164:LYS:HE3	2.02	0.41
1:B:11:THR:HB	1:B:61:LEU:HD11	2.01	0.41
1:G:32:ILE:HG22	1:G:34:LEU:HG	2.02	0.41
1:A:10:VAL:HG12	1:A:13:SER:HB3	2.03	0.41
1:D:110:LEU:HB3	1:D:111:ASN:HD22	1.85	0.41
1:F:57:HIS:HB2	1:F:73:LEU:HD12	2.03	0.41
1:B:141:HIS:HD2	3:B:282:HOH:O	2.02	0.41
1:C:223:THR:H	1:C:226:HIS:CD2	2.39	0.41
1:C:90:ILE:H	1:C:111:ASN:HD21	1.68	0.41
1:H:110:LEU:O	1:H:114:ALA:HB3	2.21	0.41
1:B:8:ALA:HB3	1:B:32:ILE:HD13	2.02	0.41
1:A:215:LYS:HB3	1:A:253:TRP:CZ2	2.56	0.41
1:A:73:LEU:HD11	1:A:77:GLU:HG3	2.03	0.41
1:E:86:ASN:HB2	1:E:136:ASN:ND2	2.33	0.41
1:H:92:HIS:CE1	3:H:338:HOH:O	2.64	0.41
1:E:7:THR:HG21	1:E:78:PHE:HB2	2.03	0.41
1:F:123:LEU:N	1:F:124:PRO:CD	2.83	0.41
1:E:110:LEU:O	1:E:114:ALA:HB3	2.21	0.41
1:B:187:THR:HB	1:B:188:PRO:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:119:THR:HG23	1:E:134:ILE:HD13	2.01	0.41
1:C:186:LEU:HD22	1:C:221:PHE:CE1	2.55	0.41
1:D:90:ILE:HD11	1:D:106:LYS:HG3	2.03	0.40
1:A:182:PRO:HA	1:A:249:VAL:O	2.21	0.40
1:E:61:LEU:HD22	1:E:67:ILE:HG12	2.03	0.40
1:G:47:ILE:HD12	1:G:52:VAL:HG12	2.02	0.40
1:G:35:ASN:HA	1:G:57:HIS:O	2.21	0.40
1:F:111:ASN:HD22	1:F:111:ASN:N	2.19	0.40
1:C:57:HIS:HD2	1:C:59:ALA:N	2.17	0.40
1:E:149:LYS:HZ1	1:E:193:GLN:HE22	1.69	0.40
1:C:86:ASN:HD22	1:C:136:ASN:HD21	1.69	0.40
1:G:60:ASP:H	1:G:66:GLN:NE2	2.19	0.40
1:G:14:THR:CG2	1:G:37:PHE:O	2.60	0.40
1:F:182:PRO:HB2	2:F:300:NAD:C5N	2.51	0.40
1:C:121:LEU:HD23	1:C:121:LEU:HA	1.93	0.40
1:B:203:ASP:HA	1:B:204:PRO:HD2	1.92	0.40
1:D:57:HIS:HE1	3:D:363:HOH:O	2.05	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)
3:E:336:HOH:O	3:G:332:HOH:O[2_555]	1.98	0.22

## 5.3 Torsion angles [i](#)

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

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	246/255 (96%)	237 (96%)	9 (4%)	0	100	100
1	B	248/255 (97%)	230 (93%)	15 (6%)	3 (1%)	16	8
1	C	239/255 (94%)	224 (94%)	13 (5%)	2 (1%)	24	15

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	253/255 (99%)	244 (96%)	8 (3%)	1 (0%)	39	32
1	E	245/255 (96%)	233 (95%)	11 (4%)	1 (0%)	39	32
1	F	234/255 (92%)	226 (97%)	8 (3%)	0	100	100
1	G	247/255 (97%)	237 (96%)	7 (3%)	3 (1%)	16	8
1	H	236/255 (92%)	220 (93%)	14 (6%)	2 (1%)	24	15
All	All	1948/2040 (96%)	1851 (95%)	85 (4%)	12 (1%)	30	21

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	39	ASP
1	H	46	GLU
1	H	47	ILE
1	G	207	ALA
1	B	196	ASP
1	B	130	ASN
1	B	197	ARG
1	G	138	ALA
1	C	131	TRP
1	D	202	GLY
1	E	206	GLN
1	G	206	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	188/189 (100%)	175 (93%)	13 (7%)	19	12
1	B	188/189 (100%)	175 (93%)	13 (7%)	19	12
1	C	181/189 (96%)	159 (88%)	22 (12%)	6	3
1	D	189/189 (100%)	174 (92%)	15 (8%)	15	9
1	E	186/189 (98%)	166 (89%)	20 (11%)	8	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	176/189 (93%)	161 (92%)	15 (8%)	13	7
1	G	187/189 (99%)	169 (90%)	18 (10%)	10	5
1	H	178/189 (94%)	168 (94%)	10 (6%)	26	19
All	All	1473/1512 (97%)	1347 (91%)	126 (9%)	13	7

All (126) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	23	GLN
1	A	25	LEU
1	A	27	ARG
1	A	53	LYS
1	A	98	GLN
1	A	106	LYS
1	A	111	ASN
1	A	192	LYS
1	A	205	LEU
1	A	212	LEU
1	A	217	PRO
1	A	230	LEU
1	B	25	LEU
1	B	52	VAL
1	B	53	LYS
1	B	73	LEU
1	B	76	ARG
1	B	106	LYS
1	B	186	LEU
1	B	192	LYS
1	B	194	ILE
1	B	195	ASP
1	B	197	ARG
1	B	205	LEU
1	B	225	GLU
1	C	15	SER
1	C	23	GLN
1	C	25	LEU
1	C	34	LEU
1	C	47	ILE
1	C	49	ARG
1	C	52	VAL

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Mol	Chain	Res	Type
1	C	76	ARG
1	C	102	GLU
1	C	110	LEU
1	C	111	ASN
1	C	189	LEU
1	C	190	VAL
1	C	191	GLN
1	C	192	LYS
1	C	193	GLN
1	C	210	ASP
1	C	211	LEU
1	C	212	LEU
1	C	214	GLU
1	C	230	LEU
1	C	232	LEU
1	D	2	THR
1	D	14	THR
1	D	15	SER
1	D	23	GLN
1	D	25	LEU
1	D	34	LEU
1	D	76	ARG
1	D	101	LEU
1	D	102	GLU
1	D	110	LEU
1	D	111	ASN
1	D	188	PRO
1	D	205	LEU
1	D	230	LEU
1	D	232	LEU
1	E	6	LYS
1	E	14	THR
1	E	25	LEU
1	E	27	ARG
1	E	42	PRO
1	E	47	ILE
1	E	49	ARG
1	E	75	GLU
1	E	98	GLN
1	E	103	SER
1	E	110	LEU
1	E	120	ARG

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Mol	Chain	Res	Type
1	E	124	PRO
1	E	130	ASN
1	E	192	LYS
1	E	206	GLN
1	E	212	LEU
1	E	225	GLU
1	E	230	LEU
1	E	232	LEU
1	F	2	THR
1	F	23	GLN
1	F	25	LEU
1	F	49	ARG
1	F	52	VAL
1	F	73	LEU
1	F	76	ARG
1	F	97	GLU
1	F	101	LEU
1	F	110	LEU
1	F	111	ASN
1	F	140	VAL
1	F	186	LEU
1	F	212	LEU
1	F	230	LEU
1	G	2	THR
1	G	13	SER
1	G	14	THR
1	G	25	LEU
1	G	47	ILE
1	G	52	VAL
1	G	96	VAL
1	G	106	LYS
1	G	110	LEU
1	G	111	ASN
1	G	192	LYS
1	G	208	GLN
1	G	210	ASP
1	G	211	LEU
1	G	212	LEU
1	G	214	GLU
1	G	230	LEU
1	G	232	LEU
1	H	23	GLN

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Mol	Chain	Res	Type
1	H	25	LEU
1	H	46	GLU
1	H	101	LEU
1	H	102	GLU
1	H	110	LEU
1	H	111	ASN
1	H	206	GLN
1	H	212	LEU
1	H	230	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (73) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	35	ASN
1	A	66	GLN
1	A	111	ASN
1	A	136	ASN
1	A	141	HIS
1	A	216	GLN
1	A	226	HIS
1	A	256	GLN
1	B	35	ASN
1	B	66	GLN
1	B	111	ASN
1	B	136	ASN
1	B	141	HIS
1	B	216	GLN
1	B	226	HIS
1	B	256	GLN
1	C	35	ASN
1	C	57	HIS
1	C	66	GLN
1	C	111	ASN
1	C	136	ASN
1	C	141	HIS
1	C	157	HIS
1	C	193	GLN
1	C	216	GLN
1	D	35	ASN
1	D	57	HIS
1	D	66	GLN

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Mol	Chain	Res	Type
1	D	92	HIS
1	D	98	GLN
1	D	111	ASN
1	D	130	ASN
1	D	136	ASN
1	D	141	HIS
1	D	157	HIS
1	D	193	GLN
1	D	216	GLN
1	E	35	ASN
1	E	57	HIS
1	E	66	GLN
1	E	98	GLN
1	E	111	ASN
1	E	136	ASN
1	E	141	HIS
1	E	157	HIS
1	E	193	GLN
1	E	216	GLN
1	E	226	HIS
1	E	256	GLN
1	F	35	ASN
1	F	66	GLN
1	F	111	ASN
1	F	136	ASN
1	F	141	HIS
1	F	216	GLN
1	G	35	ASN
1	G	57	HIS
1	G	66	GLN
1	G	111	ASN
1	G	136	ASN
1	G	141	HIS
1	G	157	HIS
1	G	193	GLN
1	G	216	GLN
1	H	35	ASN
1	H	66	GLN
1	H	92	HIS
1	H	111	ASN
1	H	130	ASN
1	H	136	ASN

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Mol	Chain	Res	Type
1	H	141	HIS
1	H	216	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

6 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	C	300	-	38,48,48	1.87	5 (13%)	47,73,73	2.39	8 (17%)
2	NAD	D	300	-	38,48,48	1.48	7 (18%)	47,73,73	2.79	12 (25%)
2	NAD	E	300	-	38,48,48	1.53	5 (13%)	47,73,73	2.60	11 (23%)
2	NAD	F	300	-	38,48,48	1.71	6 (15%)	47,73,73	2.57	8 (17%)
2	NAD	G	300	-	38,48,48	1.62	4 (10%)	47,73,73	1.73	6 (12%)
2	NAD	H	300	-	38,48,48	1.74	7 (18%)	47,73,73	2.71	7 (14%)

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	C	300	-	-	0/22/62/62	0/5/5/5
2	NAD	D	300	-	-	0/22/62/62	0/5/5/5
2	NAD	E	300	-	-	0/22/62/62	0/5/5/5
2	NAD	F	300	-	-	0/22/62/62	0/5/5/5
2	NAD	G	300	-	-	0/22/62/62	0/5/5/5
2	NAD	H	300	-	-	0/22/62/62	0/5/5/5

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	300	NAD	C2N-C3N	-3.06	1.34	1.39
2	D	300	NAD	PA-O2A	-2.10	1.46	1.54
2	C	300	NAD	O4D-C1D	2.02	1.43	1.41
2	H	300	NAD	O4D-C1D	2.11	1.43	1.41
2	F	300	NAD	O4B-C1B	2.18	1.44	1.41
2	H	300	NAD	C4A-N3A	2.23	1.38	1.35
2	H	300	NAD	O4B-C1B	2.28	1.44	1.41
2	G	300	NAD	C2A-N3A	2.32	1.36	1.32
2	F	300	NAD	C2N-C3N	2.32	1.42	1.39
2	E	300	NAD	C3N-C7N	2.39	1.54	1.50
2	D	300	NAD	C2A-N3A	2.40	1.36	1.32
2	E	300	NAD	C2A-N3A	2.63	1.36	1.32
2	C	300	NAD	O4B-C1B	2.64	1.44	1.41
2	E	300	NAD	O4D-C1D	2.67	1.44	1.41
2	D	300	NAD	O4B-C1B	2.73	1.44	1.41
2	H	300	NAD	C2N-C3N	2.77	1.43	1.39
2	D	300	NAD	O4D-C1D	2.88	1.44	1.41
2	C	300	NAD	C2A-N1A	2.89	1.39	1.33
2	H	300	NAD	C2A-N1A	2.94	1.39	1.33
2	F	300	NAD	O4D-C1D	3.03	1.45	1.41
2	E	300	NAD	C5N-C4N	3.15	1.45	1.38
2	H	300	NAD	C2A-N3A	3.39	1.38	1.32
2	F	300	NAD	C2A-N1A	3.43	1.40	1.33
2	G	300	NAD	O4B-C1B	3.59	1.45	1.41
2	D	300	NAD	O7N-C7N	3.61	1.31	1.24
2	C	300	NAD	C2A-N3A	3.66	1.38	1.32
2	F	300	NAD	C2A-N3A	3.74	1.38	1.32
2	D	300	NAD	C2A-N1A	3.84	1.41	1.33
2	G	300	NAD	C2A-N1A	4.31	1.42	1.33
2	E	300	NAD	O7N-C7N	5.64	1.36	1.24
2	G	300	NAD	O7N-C7N	5.76	1.36	1.24
2	F	300	NAD	O7N-C7N	6.36	1.37	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	300	NAD	O7N-C7N	7.13	1.39	1.24
2	C	300	NAD	O7N-C7N	8.83	1.43	1.24

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	300	NAD	N3A-C2A-N1A	-14.74	117.61	128.89
2	C	300	NAD	N3A-C2A-N1A	-11.44	120.13	128.89
2	E	300	NAD	N3A-C2A-N1A	-10.40	120.93	128.89
2	F	300	NAD	N3A-C2A-N1A	-9.98	121.25	128.89
2	D	300	NAD	N3A-C2A-N1A	-8.60	122.31	128.89
2	G	300	NAD	N3A-C2A-N1A	-6.35	124.03	128.89
2	D	300	NAD	O7N-C7N-N7N	-6.13	113.97	122.59
2	H	300	NAD	C1B-N9A-C4A	-5.53	118.60	126.94
2	E	300	NAD	O7N-C7N-C3N	-5.42	113.67	119.59
2	G	300	NAD	O7N-C7N-C3N	-5.27	113.83	119.59
2	E	300	NAD	O3-PN-O5D	-4.10	92.06	102.94
2	D	300	NAD	O7N-C7N-C3N	-4.03	115.19	119.59
2	C	300	NAD	C2B-C1B-N9A	-3.97	108.23	114.29
2	D	300	NAD	C4D-O4D-C1D	-3.68	105.67	109.72
2	D	300	NAD	O3D-C3D-C2D	-3.45	100.62	111.83
2	D	300	NAD	O3-PN-O5D	-3.42	93.86	102.94
2	C	300	NAD	O3-PN-O5D	-3.34	94.07	102.94
2	F	300	NAD	C1B-N9A-C4A	-3.29	121.98	126.94
2	C	300	NAD	PN-O3-PA	-3.13	123.94	132.73
2	F	300	NAD	O7N-C7N-N7N	-3.12	118.21	122.59
2	E	300	NAD	O7N-C7N-N7N	-2.80	118.66	122.59
2	D	300	NAD	O3B-C3B-C4B	-2.69	103.00	111.05
2	G	300	NAD	O5B-PA-O1A	-2.59	99.55	109.62
2	F	300	NAD	O5D-PN-O1N	-2.51	99.85	109.62
2	F	300	NAD	C2B-C1B-N9A	-2.49	110.49	114.29
2	E	300	NAD	O3B-C3B-C2B	-2.35	104.18	111.83
2	E	300	NAD	C1B-N9A-C4A	-2.25	123.55	126.94
2	H	300	NAD	O7N-C7N-N7N	-2.23	119.45	122.59
2	E	300	NAD	C4D-O4D-C1D	-2.18	107.32	109.72
2	D	300	NAD	C4A-C5A-N7A	-2.10	107.54	109.48
2	D	300	NAD	N6A-C6A-N1A	-2.06	114.77	119.20
2	D	300	NAD	C5N-C6N-N1N	-2.04	116.94	120.47
2	C	300	NAD	O5D-PN-O1N	2.00	117.39	109.62
2	C	300	NAD	O7N-C7N-N7N	2.01	125.42	122.59
2	D	300	NAD	O2N-PN-O1N	2.04	123.61	112.53
2	E	300	NAD	O2B-C2B-C3B	2.11	118.69	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300	NAD	O3-PA-O5B	2.11	108.54	102.94
2	G	300	NAD	C2B-C1B-N9A	2.12	117.54	114.29
2	F	300	NAD	O2B-C2B-C3B	2.17	118.90	111.83
2	E	300	NAD	C2A-N1A-C6A	2.36	122.98	118.77
2	H	300	NAD	O2A-PA-O3	2.45	116.21	105.09
2	H	300	NAD	C4D-O4D-C1D	2.50	112.47	109.72
2	G	300	NAD	O4B-C1B-N9A	2.87	114.11	108.10
2	H	300	NAD	C2A-N1A-C6A	3.32	124.70	118.77
2	G	300	NAD	C3N-C7N-N7N	4.59	122.84	117.82
2	E	300	NAD	O4D-C1D-N1N	4.63	113.22	108.13
2	H	300	NAD	C3N-C7N-N7N	4.69	122.94	117.82
2	F	300	NAD	C3N-C7N-N7N	5.82	124.18	117.82
2	C	300	NAD	O4D-C1D-N1N	7.46	116.33	108.13
2	E	300	NAD	C3N-C7N-N7N	8.78	127.43	117.82
2	F	300	NAD	O4D-C1D-N1N	9.47	118.54	108.13
2	D	300	NAD	C3N-C7N-N7N	11.81	130.74	117.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	300	NAD	1	0
2	D	300	NAD	2	0
2	E	300	NAD	4	0
2	F	300	NAD	3	0
2	G	300	NAD	2	0
2	H	300	NAD	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	250/255 (98%)	0.95	45 (18%) 2 2	36, 42, 52, 60	0
1	B	252/255 (98%)	0.78	30 (11%) 6 6	35, 42, 48, 59	0
1	C	243/255 (95%)	1.34	66 (27%) 1 1	35, 42, 50, 57	0
1	D	255/255 (100%)	0.88	48 (18%) 2 2	37, 42, 50, 59	0
1	E	249/255 (97%)	0.98	44 (17%) 2 2	36, 42, 48, 57	0
1	F	238/255 (93%)	0.94	39 (16%) 2 3	35, 42, 50, 55	0
1	G	251/255 (98%)	1.10	48 (19%) 2 2	35, 42, 49, 56	0
1	H	240/255 (94%)	0.98	48 (20%) 1 1	36, 42, 53, 63	0
All	All	1978/2040 (96%)	0.99	368 (18%) 2 2	35, 42, 50, 63	0

All (368) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	205	LEU	11.2
1	C	190	VAL	10.8
1	G	198	ALA	10.5
1	G	199	ALA	8.7
1	C	2	THR	8.6
1	C	189	LEU	8.5
1	C	194	ILE	7.8
1	G	205	LEU	7.8
1	D	201	GLY	7.8
1	E	198	ALA	7.7
1	G	204	PRO	7.6
1	D	200	ASN	7.4
1	H	188	PRO	7.2
1	E	194	ILE	7.1
1	C	38	GLY	7.0
1	A	194	ILE	6.9

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Mol	Chain	Res	Type	RSRZ
1	E	207	ALA	6.5
1	H	205	LEU	6.5
1	F	159	VAL	6.2
1	G	206	GLN	6.1
1	B	194	ILE	5.9
1	C	41	ALA	5.7
1	G	194	ILE	5.7
1	B	198	ALA	5.6
1	D	198	ALA	5.5
1	B	203	ASP	5.3
1	C	193	GLN	5.2
1	H	206	GLN	5.1
1	H	204	PRO	5.1
1	C	15	SER	5.1
1	A	195	ASP	5.0
1	D	199	ALA	5.0
1	C	76	ARG	4.9
1	F	189	LEU	4.9
1	D	137	ILE	4.9
1	B	207	ALA	4.9
1	C	188	PRO	4.9
1	A	196	ASP	4.9
1	E	159	VAL	4.8
1	A	52	VAL	4.7
1	H	159	VAL	4.7
1	B	202	GLY	4.6
1	B	204	PRO	4.6
1	H	4	LYS	4.6
1	B	197	ARG	4.6
1	H	52	VAL	4.6
1	D	159	VAL	4.5
1	D	160	VAL	4.5
1	G	159	VAL	4.5
1	F	211	LEU	4.4
1	F	208	GLN	4.4
1	A	54	ALA	4.4
1	B	205	LEU	4.4
1	E	160	VAL	4.4
1	H	2	THR	4.4
1	G	160	VAL	4.3
1	E	76	ARG	4.3
1	F	188	PRO	4.3

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Mol	Chain	Res	Type	RSRZ
1	E	196	ASP	4.3
1	A	205	LEU	4.3
1	C	209	HIS	4.3
1	A	206	GLN	4.2
1	C	208	GLN	4.2
1	G	2	THR	4.2
1	C	3	LEU	4.2
1	F	186	LEU	4.2
1	D	180	ILE	4.1
1	B	189	LEU	4.1
1	A	197	ARG	4.0
1	C	186	LEU	4.1
1	G	135	ILE	4.0
1	G	76	ARG	4.0
1	C	55	VAL	4.0
1	D	162	LEU	4.0
1	H	76	ARG	4.0
1	B	190	VAL	4.0
1	E	206	GLN	4.0
1	E	155	ALA	4.0
1	A	204	PRO	3.9
1	C	211	LEU	3.9
1	D	135	ILE	3.8
1	F	115	VAL	3.8
1	G	137	ILE	3.8
1	F	160	VAL	3.8
1	D	112	LEU	3.8
1	A	65	ALA	3.7
1	C	42	PRO	3.7
1	E	154	ALA	3.7
1	G	207	ALA	3.7
1	C	48	ALA	3.7
1	D	181	CYS	3.7
1	F	112	LEU	3.7
1	A	203	ASP	3.7
1	D	84	LEU	3.7
1	D	179	ALA	3.7
1	H	155	ALA	3.7
1	H	160	VAL	3.7
1	D	138	ALA	3.7
1	F	2	THR	3.6
1	C	137	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	159	VAL	3.6
1	A	49	ARG	3.6
1	E	5	GLY	3.5
1	D	85	VAL	3.5
1	E	162	LEU	3.5
1	G	163	THR	3.5
1	G	196	ASP	3.5
1	C	4	LYS	3.5
1	E	143	LEU	3.5
1	B	160	VAL	3.4
1	A	211	LEU	3.4
1	D	163	THR	3.4
1	F	137	ILE	3.4
1	A	190	VAL	3.4
1	C	81	VAL	3.4
1	G	55	VAL	3.4
1	G	143	LEU	3.4
1	H	207	ALA	3.4
1	E	137	ILE	3.4
1	A	189	LEU	3.4
1	D	115	VAL	3.3
1	E	197	ARG	3.3
1	G	192	LYS	3.3
1	F	209	HIS	3.3
1	F	162	LEU	3.3
1	H	112	LEU	3.3
1	C	39	ASP	3.3
1	A	187	THR	3.3
1	F	158	GLY	3.3
1	H	144	VAL	3.3
1	A	46	GLU	3.3
1	A	192	LYS	3.2
1	H	80	GLY	3.2
1	C	128	ALA	3.2
1	H	208	GLN	3.2
1	G	162	LEU	3.2
1	D	155	ALA	3.2
1	A	51	GLY	3.2
1	H	158	GLY	3.2
1	E	209	HIS	3.2
1	G	155	ALA	3.2
1	D	136	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	202	GLY	3.2
1	A	207	ALA	3.2
1	G	195	ASP	3.2
1	C	192	LYS	3.2
1	F	210	ASP	3.2
1	G	177	CYS	3.2
1	E	208	GLN	3.2
1	G	65	ALA	3.1
1	H	51	GLY	3.1
1	B	196	ASP	3.1
1	C	73	LEU	3.1
1	C	80	GLY	3.1
1	F	155	ALA	3.1
1	F	163	THR	3.1
1	C	54	ALA	3.1
1	A	2	THR	3.1
1	F	135	ILE	3.1
1	G	134	ILE	3.1
1	E	190	VAL	3.1
1	H	48	ALA	3.0
1	A	208	GLN	3.0
1	C	72	ALA	3.0
1	D	246	ALA	3.0
1	E	153	VAL	3.0
1	B	85	VAL	3.0
1	E	144	VAL	3.0
1	B	143	LEU	3.0
1	D	205	LEU	3.0
1	D	144	VAL	3.0
1	A	45	ALA	3.0
1	C	160	VAL	2.9
1	G	190	VAL	2.9
1	D	247	TRP	2.9
1	D	140	VAL	2.9
1	E	158	GLY	2.9
1	F	38	GLY	2.9
1	H	41	ALA	2.9
1	A	42	PRO	2.9
1	C	64	VAL	2.9
1	C	138	ALA	2.9
1	B	192	LYS	2.9
1	G	53	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	85	VAL	2.9
1	A	160	VAL	2.9
1	F	85	VAL	2.9
1	E	138	ALA	2.9
1	H	154	ALA	2.8
1	G	197	ARG	2.8
1	H	142	GLY	2.8
1	D	204	PRO	2.8
1	E	192	LYS	2.8
1	H	45	ALA	2.8
1	A	158	GLY	2.8
1	C	187	THR	2.8
1	B	74	ALA	2.8
1	D	142	GLY	2.8
1	E	195	ASP	2.8
1	G	85	VAL	2.8
1	E	210	ASP	2.8
1	G	208	GLN	2.7
1	H	39	ASP	2.7
1	C	47	ILE	2.7
1	G	179	ALA	2.7
1	H	186	LEU	2.7
1	C	52	VAL	2.7
1	C	37	PHE	2.7
1	B	155	ALA	2.7
1	E	179	ALA	2.7
1	H	138	ALA	2.7
1	A	209	HIS	2.7
1	D	87	ASN	2.7
1	F	49	ARG	2.7
1	H	161	GLY	2.6
1	B	195	ASP	2.6
1	C	69	ALA	2.6
1	C	207	ALA	2.6
1	A	80	GLY	2.6
1	F	161	GLY	2.6
1	G	77	GLU	2.6
1	E	115	VAL	2.6
1	G	166	VAL	2.6
1	G	193	GLN	2.6
1	C	49	ARG	2.6
1	C	5	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	50	HIS	2.6
1	H	115	VAL	2.6
1	D	86	ASN	2.5
1	D	242	VAL	2.5
1	E	135	ILE	2.5
1	E	157	HIS	2.5
1	H	209	HIS	2.5
1	C	63	ASP	2.5
1	F	180	ILE	2.5
1	H	143	LEU	2.5
1	D	156	LYS	2.5
1	D	9	LEU	2.5
1	A	38	GLY	2.5
1	G	144	VAL	2.5
1	G	180	ILE	2.5
1	E	163	THR	2.5
1	D	207	ALA	2.5
1	C	101	LEU	2.5
1	G	39	ASP	2.5
1	D	157	HIS	2.5
1	D	153	VAL	2.5
1	G	115	VAL	2.5
1	C	191	GLN	2.5
1	F	177	CYS	2.5
1	D	130	ASN	2.4
1	C	43	ALA	2.4
1	D	177	CYS	2.4
1	F	245	ALA	2.4
1	C	158	GLY	2.4
1	G	161	GLY	2.4
1	H	157	HIS	2.4
1	G	210	ASP	2.4
1	H	187	THR	2.4
1	C	67	ILE	2.4
1	A	191	GLN	2.4
1	B	76	ARG	2.4
1	G	191	GLN	2.4
1	E	246	ALA	2.4
1	F	154	ALA	2.4
1	F	136	ASN	2.4
1	F	179	ALA	2.4
1	H	86	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	37	PHE	2.4
1	A	186	LEU	2.4
1	A	48	ALA	2.3
1	C	58	PRO	2.3
1	F	96	VAL	2.3
1	H	137	ILE	2.3
1	B	73	LEU	2.3
1	C	210	ASP	2.3
1	C	45	ALA	2.3
1	H	114	ALA	2.3
1	C	87	ASN	2.3
1	C	92	HIS	2.3
1	H	163	THR	2.3
1	A	76	ARG	2.3
1	E	181	CYS	2.3
1	C	85	VAL	2.3
1	E	140	VAL	2.3
1	H	55	VAL	2.3
1	E	85	VAL	2.3
1	H	11	THR	2.3
1	C	28	ALA	2.3
1	C	155	ALA	2.3
1	F	63	ASP	2.3
1	B	80	GLY	2.3
1	C	102	GLU	2.2
1	D	166	VAL	2.2
1	C	74	ALA	2.2
1	E	72	ALA	2.2
1	E	161	GLY	2.2
1	G	74	ALA	2.2
1	B	112	LEU	2.2
1	H	162	LEU	2.2
1	C	53	LYS	2.2
1	D	158	GLY	2.2
1	F	214	GLU	2.2
1	G	41	ALA	2.2
1	A	53	LYS	2.2
1	C	65	ALA	2.2
1	H	151	ALA	2.2
1	D	203	ASP	2.2
1	D	143	LEU	2.2
1	E	84	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	152	TYR	2.2
1	A	161	GLY	2.2
1	G	79	GLY	2.2
1	G	244	GLY	2.2
1	A	193	GLN	2.2
1	C	10	VAL	2.2
1	B	245	ALA	2.2
1	F	207	ALA	2.2
1	G	246	ALA	2.2
1	H	254	LEU	2.2
1	C	163	THR	2.2
1	E	2	THR	2.2
1	C	159	VAL	2.2
1	F	138	ALA	2.2
1	H	179	ALA	2.2
1	F	134	ILE	2.2
1	C	232	LEU	2.1
1	C	77	GLU	2.1
1	A	153	VAL	2.1
1	A	159	VAL	2.1
1	E	255	ALA	2.1
1	H	150	ALA	2.1
1	H	58	PRO	2.1
1	B	180	ILE	2.1
1	B	206	GLN	2.1
1	A	74	ALA	2.1
1	E	77	GLU	2.1
1	D	182	PRO	2.1
1	H	153	VAL	2.1
1	B	162	LEU	2.1
1	G	211	LEU	2.1
1	C	8	ALA	2.1
1	H	140	VAL	2.1
1	F	11	THR	2.1
1	D	245	ALA	2.1
1	C	97	GLU	2.1
1	F	10	VAL	2.1
1	B	10	VAL	2.1
1	B	247	TRP	2.0
1	F	77	GLU	2.0
1	E	180	ILE	2.0
1	A	58	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	84	LEU	2.0
1	F	84	LEU	2.0
1	E	80	GLY	2.0
1	G	45	ALA	2.0
1	D	102	GLU	2.0
1	G	102	GLU	2.0
1	A	47	ILE	2.0
1	H	180	ILE	2.0
1	C	121	LEU	2.0
1	E	219	LEU	2.0
1	A	39	ASP	2.0
1	D	165	VAL	2.0
1	F	140	VAL	2.0
1	C	51	GLY	2.0
1	D	17	ILE	2.0
1	D	195	ASP	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAD	C	300	44/44	0.80	0.16	-0.44	47,62,73,77	0
2	NAD	F	300	44/44	0.90	0.14	-0.51	39,51,56,59	0
2	NAD	D	300	44/44	0.95	0.11	-0.78	25,35,41,44	0
2	NAD	G	300	44/44	0.95	0.09	-1.04	35,42,48,52	0
2	NAD	E	300	44/44	0.95	0.09	-1.11	30,39,45,49	0
2	NAD	H	300	44/44	0.94	0.09	-1.12	33,42,47,51	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.