



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:25 AM GMT

PDB ID : 2GVG  
Title : Crystal Structure of human NMPRTase and its complex with NMN  
Authors : Khan, J.A.; Tao, X.; Tong, L.  
Deposited on : 2006-05-02  
Resolution : 2.20 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.

---

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

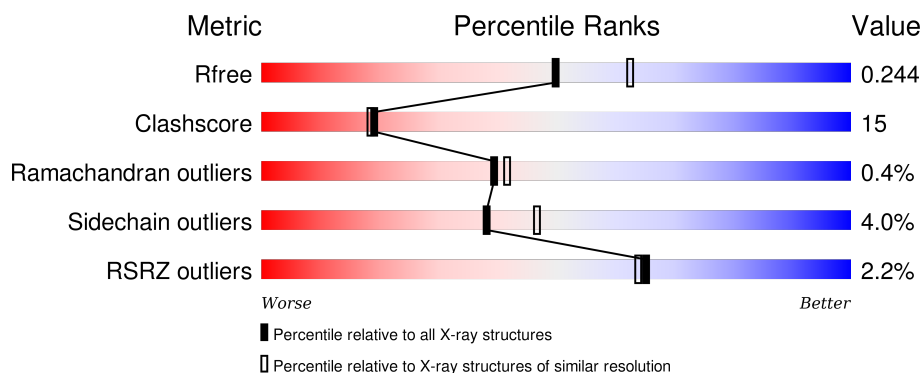
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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	491	
1	B	491	
1	C	491	
1	D	491	
1	E	491	

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	F	491	 % 71% 22% • 6%

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	PO4	E	608	-	-	-	X
3	NMN	E	505	-	-	-	X

## 2 Entry composition [i](#)

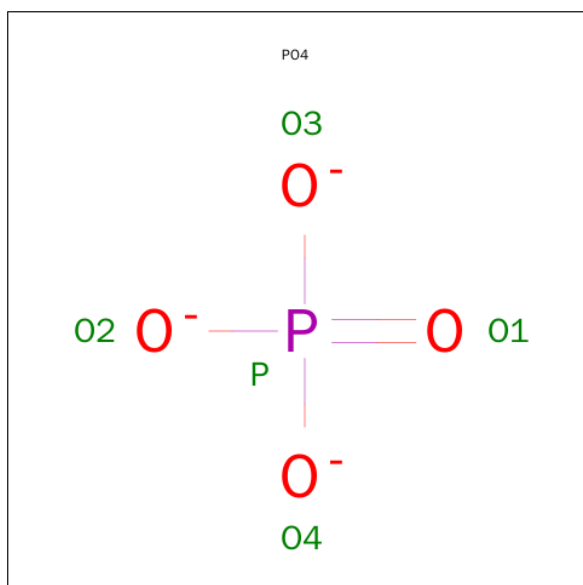
There are 4 unique types of molecules in this entry. The entry contains 23801 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nicotinamide phosphoribosyltransferase.

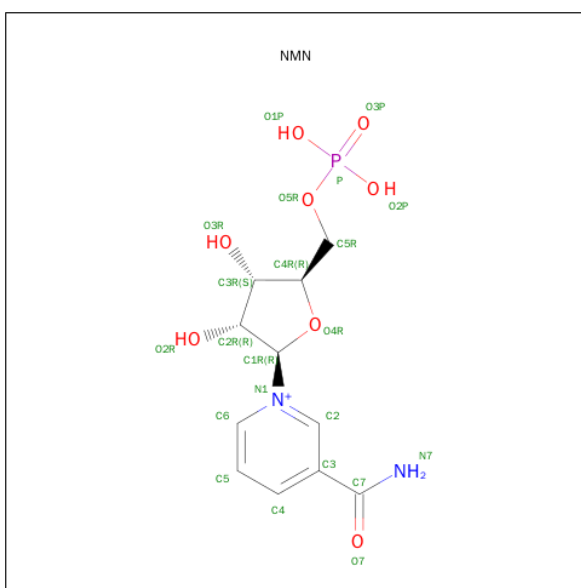
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	0	0
			3688	2376	609	696	7			
1	B	464	Total	C	N	O	S	0	0	0
			3711	2388	613	703	7			
1	C	461	Total	C	N	O	S	0	0	0
			3688	2376	609	696	7			
1	D	464	Total	C	N	O	S	0	0	0
			3711	2388	613	703	7			
1	E	464	Total	C	N	O	S	0	0	0
			3711	2388	613	703	7			
1	F	463	Total	C	N	O	S	0	0	0
			3705	2385	612	701	7			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	E	1	Total O P 5 4 1	0	0
2	D	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0

- Molecule 3 is BETA-NICOTINAMIDE RIBOSE MONOPHOSPHATE (three-letter code: NMN) (formula:  $C_{11}H_{16}N_2O_8P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
3	B	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
3	C	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
3	D	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
3	E	1	Total	C	N	O	P	0	0
			22	11	2	8	1		
3	F	1	Total	C	N	O	P	0	0
			22	11	2	8	1		

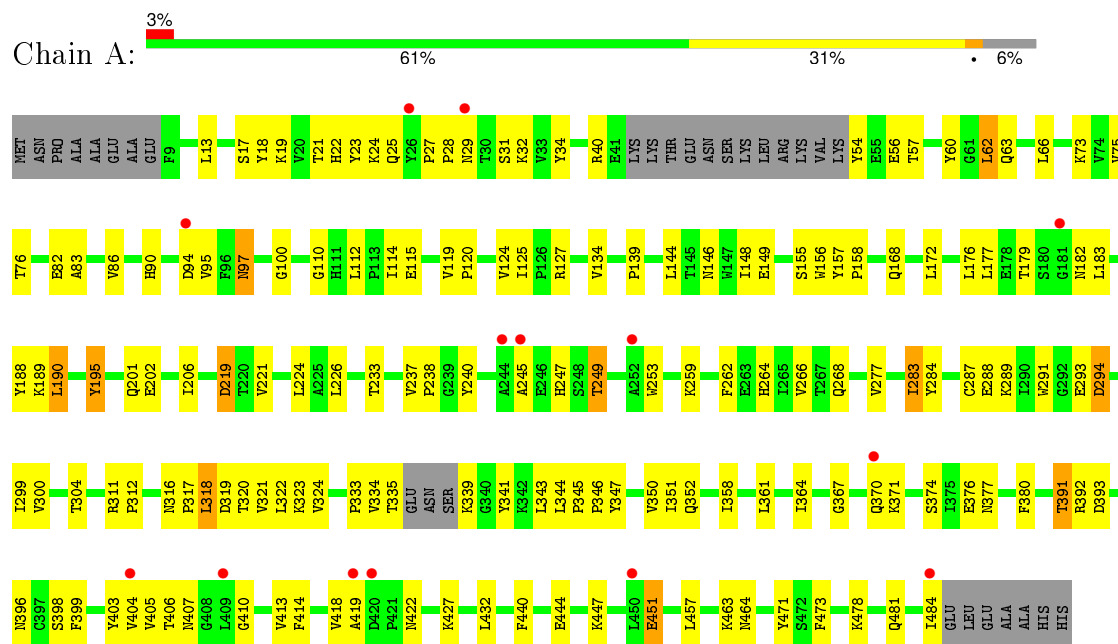
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	139	Total	O	0	0
			139	139		
4	B	165	Total	O	0	0
			165	165		
4	C	125	Total	O	0	0
			125	125		
4	D	306	Total	O	0	0
			306	306		
4	E	280	Total	O	0	0
			280	280		
4	F	380	Total	O	0	0
			380	380		

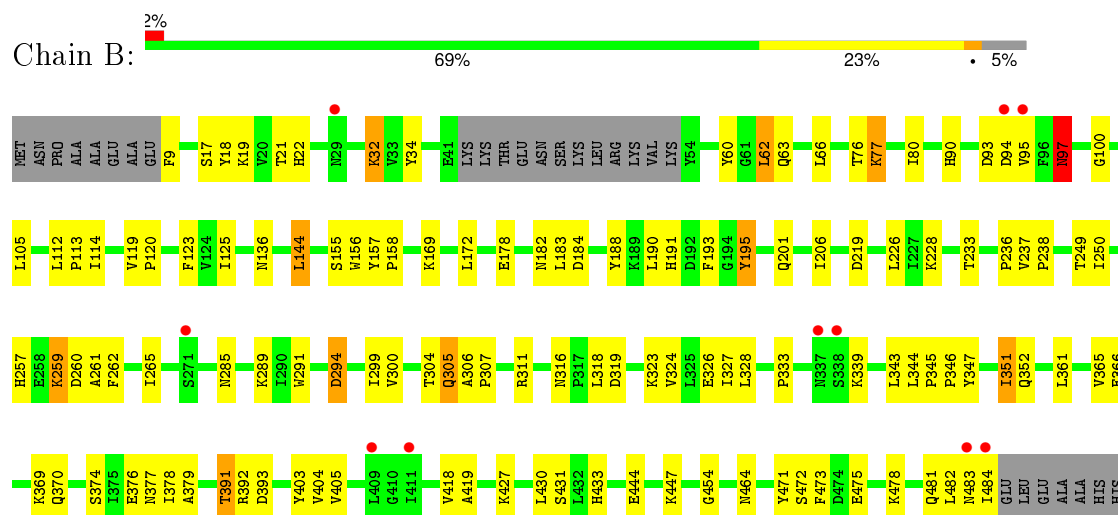
### 3 Residue-property plots

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.

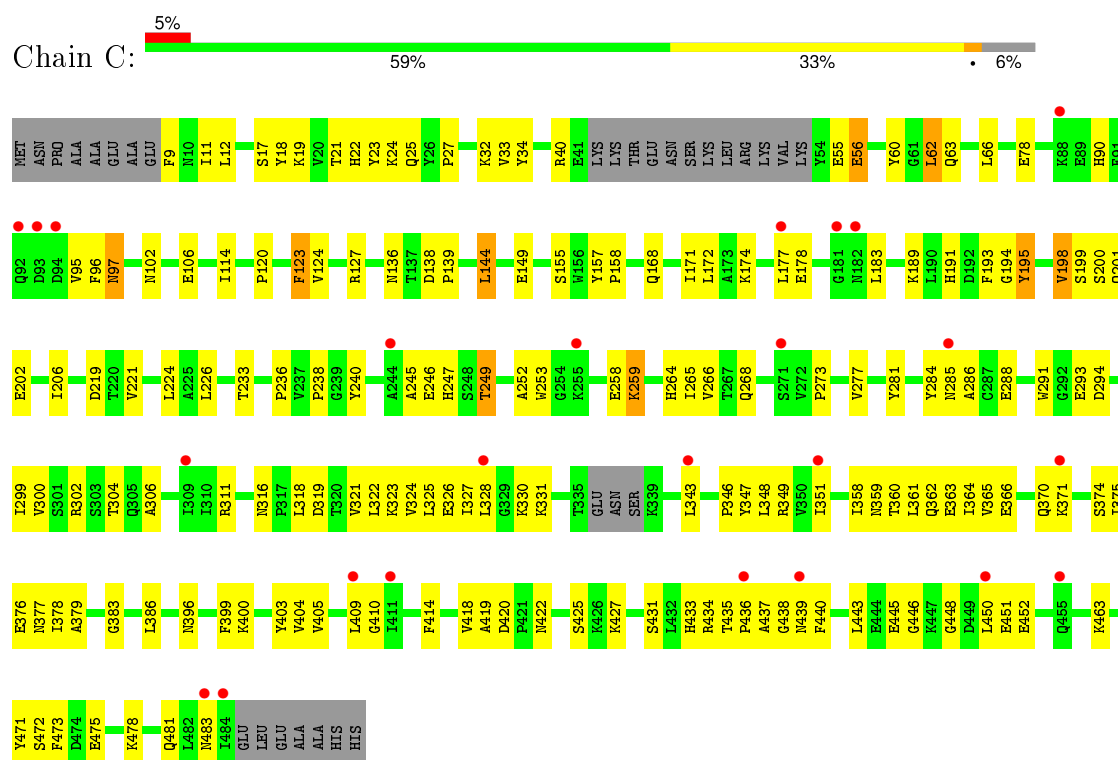
- Molecule 1: Nicotinamide phosphoribosyltransferase



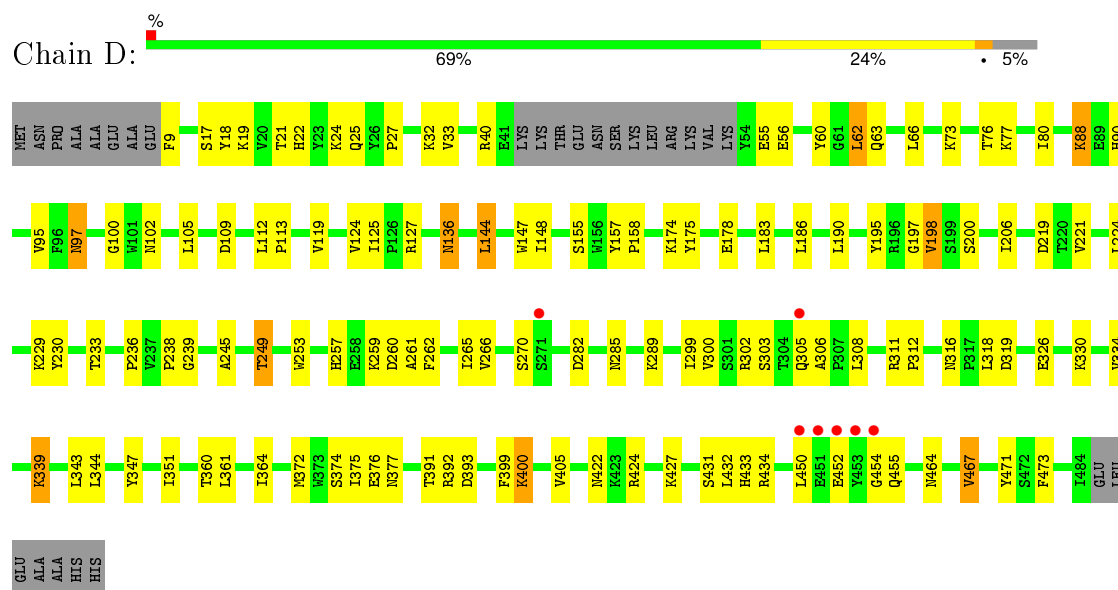
- Molecule 1: Nicotinamide phosphoribosyltransferase



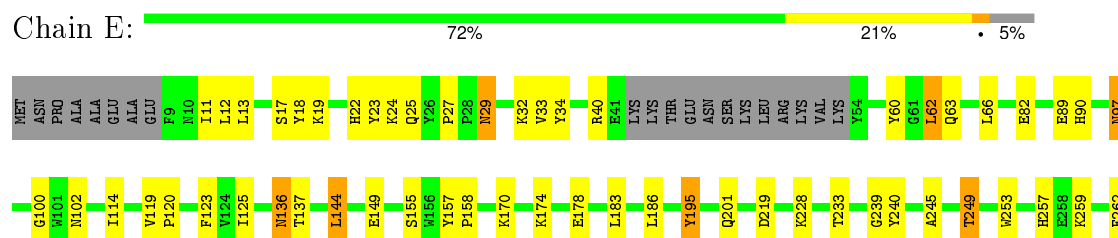
- Molecule 1: Nicotinamide phosphoribosyltransferase



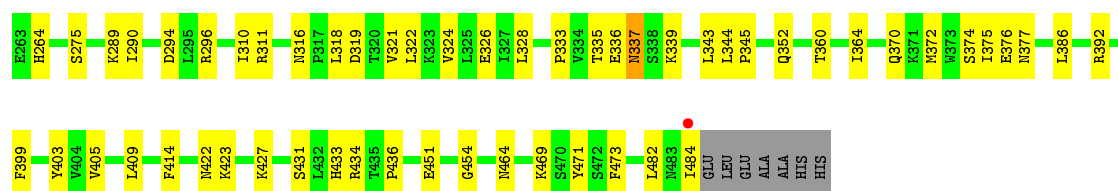
• Molecule 1: Nicotinamide phosphoribosyltransferase



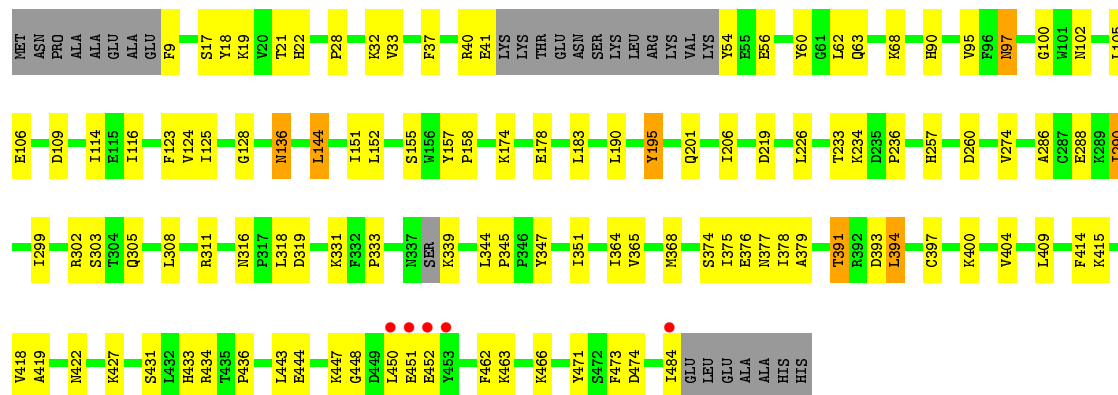
• Molecule 1: Nicotinamide phosphoribosyltransferase







• Molecule 1: Nicotinamide phosphoribosyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	253.07Å 101.37Å 148.20Å 90.00° 125.48° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 28.91 – 2.19	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.20) 92.3 (28.91-2.19)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.71 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.197 , 0.246 0.198 , 0.244	Depositor DCC
$R_{free}$ test set	10778 reflections (7.49%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.4	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 51.4	EDS
Estimated twinning fraction	0.015 for -h-2*l,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 150406 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	23801	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	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 39.91 % 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 2.9568e-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 [i](#)

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

Bond lengths and bond angles in the following residue types are not validated in this section: NMN, PO4

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	0.32	0/3775	0.57	0/5115
1	B	0.32	0/3799	0.59	0/5149
1	C	0.31	0/3775	0.56	0/5115
1	D	0.37	0/3799	0.63	1/5149 (0.0%)
1	E	0.37	0/3799	0.64	0/5149
1	F	0.39	0/3792	0.64	0/5138
All	All	0.35	0/22739	0.61	1/30815 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	127	ARG	NE-CZ-NH1	-5.51	117.55	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3688	0	3668	149	0
1	B	3711	0	3686	111	0
1	C	3688	0	3668	159	0
1	D	3711	0	3686	111	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3711	0	3686	98	0
1	F	3705	0	3680	93	0
2	A	10	0	0	1	0
2	B	10	0	0	0	0
2	C	10	0	0	1	0
2	D	10	0	0	1	0
2	E	10	0	0	0	0
2	F	10	0	0	0	0
3	A	22	0	12	0	0
3	B	22	0	12	0	0
3	C	22	0	12	1	0
3	D	22	0	12	1	0
3	E	22	0	12	1	0
3	F	22	0	12	1	0
4	A	139	0	0	9	0
4	B	165	0	0	8	0
4	C	125	0	0	10	0
4	D	306	0	0	14	0
4	E	280	0	0	9	0
4	F	380	0	0	14	0
All	All	23801	0	22146	678	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 15.

All (678) 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:D:391:THR:HG22	1:D:393:ASP:H	1.19	1.08
1:A:63:GLN:HE22	1:A:471:TYR:H	1.03	0.99
1:D:25:GLN:HE21	1:E:249:THR:HG21	1.31	0.95
1:E:344:LEU:H	1:E:377:ASN:HD21	1.08	0.95
1:E:119:VAL:HG21	1:E:125:ILE:HD11	1.46	0.93
1:B:391:THR:HG23	1:B:393:ASP:H	1.31	0.93
1:F:344:LEU:H	1:F:377:ASN:HD21	1.05	0.93
1:B:63:GLN:HE22	1:B:471:TYR:H	1.12	0.92
1:A:344:LEU:H	1:A:377:ASN:HD21	1.18	0.91
1:B:119:VAL:HG21	1:B:125:ILE:HD11	1.50	0.91
1:F:63:GLN:HE22	1:F:471:TYR:H	1.14	0.91
1:D:374:SER:OG	1:D:376:GLU:HG2	1.71	0.91
1:D:249:THR:HG21	1:E:25:GLN:NE2	1.86	0.91

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:GLN:HE22	1:C:471:TYR:H	1.19	0.90
1:B:344:LEU:H	1:B:377:ASN:ND2	1.70	0.89
1:D:148:ILE:HG23	4:D:1329:HOH:O	1.72	0.89
1:D:261:ALA:O	1:D:265:ILE:HG12	1.74	0.88
1:F:400:LYS:HE3	4:F:1363:HOH:O	1.74	0.88
1:D:119:VAL:HG21	1:D:125:ILE:HD11	1.56	0.88
1:F:374:SER:OG	1:F:376:GLU:HG2	1.74	0.88
1:A:316:ASN:HD22	1:A:319:ASP:H	1.19	0.87
1:E:316:ASN:HD22	1:E:319:ASP:H	1.22	0.87
1:A:224:LEU:HD12	1:A:238:PRO:HD2	1.57	0.87
1:D:316:ASN:HD22	1:D:319:ASP:H	1.14	0.86
1:D:25:GLN:NE2	1:E:249:THR:HG21	1.91	0.86
1:E:344:LEU:H	1:E:377:ASN:ND2	1.75	0.85
1:F:316:ASN:HD22	1:F:319:ASP:H	1.24	0.85
1:B:344:LEU:H	1:B:377:ASN:HD21	1.22	0.85
1:D:344:LEU:H	1:D:377:ASN:HD21	1.23	0.85
1:A:25:GLN:HE21	1:C:249:THR:HG21	1.40	0.84
1:D:249:THR:HG21	1:E:25:GLN:HE21	1.43	0.84
1:B:316:ASN:HD22	1:B:319:ASP:H	1.25	0.83
1:E:63:GLN:HE22	1:E:471:TYR:H	1.25	0.83
1:E:335:THR:HG23	1:E:343:LEU:HB3	1.59	0.83
1:D:63:GLN:HE22	1:D:471:TYR:H	1.22	0.82
1:C:343:LEU:HD12	1:C:377:ASN:HD21	1.41	0.81
1:A:391:THR:HG23	1:A:393:ASP:H	1.45	0.81
1:A:344:LEU:H	1:A:377:ASN:ND2	1.77	0.81
1:A:25:GLN:HE22	1:C:245:ALA:HA	1.44	0.81
1:A:249:THR:HG21	1:C:25:GLN:NE2	1.95	0.80
1:A:432:LEU:HG	1:A:457:LEU:HD12	1.63	0.79
1:D:344:LEU:H	1:D:377:ASN:ND2	1.80	0.79
1:F:56:GLU:HG2	1:F:124:VAL:HG12	1.64	0.79
1:C:343:LEU:HD12	1:C:377:ASN:ND2	1.99	0.78
1:C:400:LYS:HD3	1:C:425:SER:HB3	1.65	0.77
1:C:316:ASN:HD22	1:C:319:ASP:H	1.30	0.76
1:A:245:ALA:HA	1:C:25:GLN:HE22	1.49	0.76
1:D:56:GLU:HG2	1:D:124:VAL:CG1	2.17	0.75
1:C:478:LYS:HD2	1:C:481:GLN:NE2	2.01	0.75
1:D:303:SER:OG	1:D:305:GLN:HG2	1.87	0.75
1:B:157:TYR:HB3	1:B:158:PRO:HD3	1.69	0.75
1:E:321:VAL:HG23	1:E:352:GLN:HE21	1.51	0.75
1:B:259:LYS:HB2	1:B:259:LYS:HZ2	1.50	0.74
1:D:316:ASN:ND2	1:D:319:ASP:H	1.84	0.74

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:GLN:NE2	1:C:249:THR:HG21	2.02	0.74
1:C:478:LYS:HA	1:C:481:GLN:HE21	1.51	0.74
1:D:147:TRP:HB3	4:D:1329:HOH:O	1.85	0.74
1:D:113:PRO:HA	1:D:464:ASN:HD22	1.52	0.74
1:F:344:LEU:H	1:F:377:ASN:ND2	1.84	0.74
1:B:77:LYS:HD2	1:B:105:LEU:HD21	1.69	0.73
1:A:63:GLN:NE2	1:A:471:TYR:H	1.83	0.73
1:C:224:LEU:HD12	1:C:238:PRO:HD2	1.70	0.73
1:A:34:TYR:HB3	1:A:403:TYR:HB3	1.70	0.73
1:B:32:LYS:O	1:B:404:VAL:HG23	1.89	0.73
1:C:322:LEU:HD22	1:C:371:LYS:HG3	1.70	0.72
1:D:25:GLN:HE22	1:E:245:ALA:HA	1.54	0.72
1:F:316:ASN:ND2	1:F:319:ASP:H	1.85	0.72
1:A:63:GLN:HE22	1:A:471:TYR:N	1.86	0.72
1:E:316:ASN:ND2	1:E:319:ASP:H	1.87	0.72
1:B:259:LYS:HB2	1:B:259:LYS:NZ	2.05	0.72
1:A:344:LEU:N	1:A:377:ASN:HD21	1.87	0.71
1:E:434:ARG:HD2	4:E:1827:HOH:O	1.90	0.71
1:D:245:ALA:HA	1:E:25:GLN:HE22	1.56	0.70
1:F:183:LEU:HD11	1:F:484:ILE:HG22	1.73	0.70
1:E:195:TYR:OH	1:E:201:GLN:NE2	2.23	0.70
1:A:54:TYR:O	1:A:127:ARG:HD2	1.91	0.70
1:F:123:PHE:HB3	1:F:125:ILE:HD11	1.73	0.70
1:C:277:VAL:HB	1:C:311:ARG:NH1	2.08	0.69
1:B:374:SER:OG	1:B:376:GLU:HG2	1.91	0.69
1:F:157:TYR:HB3	1:F:158:PRO:HD3	1.75	0.69
1:E:374:SER:OG	1:E:376:GLU:HG2	1.91	0.69
1:F:257:HIS:HD2	1:F:260:ASP:OD2	1.76	0.69
1:E:157:TYR:HB3	1:E:158:PRO:HD3	1.74	0.68
1:E:316:ASN:HB3	1:E:319:ASP:HB3	1.76	0.67
1:D:56:GLU:HG2	1:D:124:VAL:HG12	1.74	0.67
1:F:286:ALA:HA	1:F:290:ILE:HG13	1.76	0.67
1:F:233:THR:HG22	1:F:473:PHE:HB3	1.75	0.67
1:D:174:LYS:O	1:D:178:GLU:HG3	1.94	0.67
1:A:157:TYR:HB3	1:A:158:PRO:HD3	1.77	0.67
1:D:424:ARG:HD2	4:D:1835:HOH:O	1.94	0.67
1:C:17:SER:OG	1:C:90:HIS:HE1	1.79	0.66
1:B:250:ILE:HD13	1:B:262:PHE:HE1	1.60	0.66
1:C:291:TRP:CE3	1:C:299:ILE:HD11	2.31	0.66
1:C:478:LYS:HD2	1:C:481:GLN:HE22	1.61	0.66
1:E:62:LEU:HD22	1:E:66:LEU:HG	1.77	0.66

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:56:GLU:HG2	1:F:124:VAL:CG1	2.25	0.66
1:F:174:LYS:O	1:F:178:GLU:HG3	1.96	0.66
1:B:391:THR:HG23	1:B:393:ASP:N	2.07	0.65
1:A:156:TRP:CG	1:A:392:ARG:HG3	2.32	0.65
1:A:249:THR:HG21	1:C:25:GLN:HE21	1.60	0.65
1:B:32:LYS:HA	1:B:136:ASN:HD21	1.61	0.65
1:E:337:ASN:C	1:E:337:ASN:HD22	2.00	0.65
1:A:21:THR:HG22	1:A:95:VAL:HG11	1.78	0.65
1:D:233:THR:HG22	1:D:473:PHE:HB3	1.79	0.65
1:E:344:LEU:N	1:E:377:ASN:HD21	1.89	0.65
1:A:291:TRP:CE3	1:A:299:ILE:HD11	2.32	0.65
1:A:195:TYR:OH	1:A:201:GLN:NE2	2.30	0.65
1:D:343:LEU:HD13	1:D:376:GLU:HG3	1.79	0.64
1:A:40:ARG:HD3	1:A:422:ASN:O	1.98	0.64
1:C:195:TYR:OH	1:C:201:GLN:NE2	2.30	0.64
1:D:249:THR:CG2	1:E:25:GLN:HE21	2.09	0.64
1:C:259:LYS:HB2	1:C:259:LYS:NZ	2.12	0.64
1:E:29:ASN:N	1:E:29:ASN:HD22	1.95	0.64
1:C:233:THR:HG22	1:C:473:PHE:HB3	1.79	0.64
1:B:119:VAL:HG21	1:B:125:ILE:CD1	2.26	0.64
1:C:114:ILE:HG12	1:C:144:LEU:HD13	1.80	0.64
1:F:391:THR:CG2	1:F:393:ASP:H	2.10	0.64
1:A:27:PRO:HG3	1:C:253:TRP:CD1	2.32	0.64
1:C:221:VAL:HA	1:C:224:LEU:HD23	1.78	0.63
1:F:63:GLN:HE22	1:F:471:TYR:N	1.92	0.63
1:C:21:THR:HG22	1:C:95:VAL:HG11	1.80	0.63
1:F:339:LYS:N	4:F:1790:HOH:O	2.30	0.63
1:C:32:LYS:HA	1:C:136:ASN:HD21	1.62	0.63
1:A:374:SER:OG	1:A:376:GLU:HG2	1.97	0.63
1:B:306:ALA:N	1:B:307:PRO:HD3	2.13	0.63
1:A:322:LEU:HD22	1:A:371:LYS:HG3	1.80	0.63
1:D:25:GLN:HE21	1:E:249:THR:CG2	2.10	0.63
1:E:155:SER:O	1:E:158:PRO:HD2	1.99	0.63
1:B:447:LYS:HG2	4:B:1537:HOH:O	1.99	0.63
1:C:19:LYS:HA	1:C:22:HIS:CD2	2.34	0.63
1:E:40:ARG:HD3	1:E:422:ASN:O	1.98	0.62
1:D:157:TYR:HB3	1:D:158:PRO:HD3	1.81	0.62
1:C:23:TYR:CE2	1:C:97:ASN:HB2	2.33	0.62
1:C:304:THR:HG23	1:C:346:PRO:HB2	1.81	0.62
1:F:311:ARG:NH2	3:F:506:NMN:O2R	2.32	0.62
1:B:195:TYR:OH	1:B:201:GLN:NE2	2.32	0.62

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:GLN:HB3	4:B:1625:HOH:O	1.99	0.62
1:E:136:ASN:HD22	1:E:136:ASN:H	1.47	0.62
1:C:157:TYR:HB3	1:C:158:PRO:HD3	1.81	0.62
1:A:316:ASN:ND2	1:A:319:ASP:H	1.95	0.62
1:F:195:TYR:OH	1:F:201:GLN:NE2	2.33	0.62
1:A:240:TYR:HA	1:C:90:HIS:CD2	2.35	0.62
1:E:32:LYS:HB3	1:E:405:VAL:HB	1.81	0.62
1:B:344:LEU:N	1:B:377:ASN:HD21	1.96	0.61
1:D:33:VAL:H	1:D:136:ASN:ND2	1.98	0.61
1:B:169:LYS:HG2	1:B:482:LEU:HD11	1.82	0.61
1:A:90:HIS:CD2	1:C:240:TYR:HA	2.35	0.61
1:E:114:ILE:HG12	1:E:144:LEU:HD13	1.83	0.61
1:B:93:ASP:OD1	1:B:94:ASP:N	2.34	0.61
1:A:112:LEU:O	1:A:464:ASN:HA	2.01	0.60
1:C:63:GLN:HE22	1:C:471:TYR:N	1.96	0.60
1:B:114:ILE:HD13	1:B:136:ASN:HA	1.83	0.60
1:A:119:VAL:HG21	1:A:125:ILE:HD11	1.83	0.60
1:D:391:THR:HG22	1:D:393:ASP:N	2.04	0.60
1:C:400:LYS:CD	1:C:425:SER:HB3	2.30	0.60
1:C:127:ARG:HH21	1:C:396:ASN:HD22	1.47	0.60
1:B:172:LEU:HD21	1:B:361:LEU:HD11	1.84	0.60
1:D:299:ILE:HD12	1:D:308:LEU:HD22	1.84	0.60
1:A:259:LYS:HE3	1:A:294:ASP:HB3	1.83	0.60
1:C:102:ASN:O	1:C:106:GLU:HG3	2.01	0.60
1:E:63:GLN:HE22	1:E:471:TYR:N	1.97	0.60
1:B:114:ILE:HG12	1:B:144:LEU:HD13	1.84	0.60
1:C:24:LYS:NZ	1:C:95:VAL:HG13	2.16	0.60
1:A:343:LEU:HD13	1:A:376:GLU:HG3	1.84	0.60
1:D:33:VAL:H	1:D:136:ASN:HD21	1.48	0.59
1:C:414:PHE:CD2	1:C:427:LYS:HG2	2.38	0.59
1:B:261:ALA:O	1:B:265:ILE:HG12	2.02	0.59
1:A:259:LYS:CE	1:A:294:ASP:HB3	2.32	0.59
1:B:431:SER:OG	1:B:433:HIS:HE1	1.86	0.59
1:D:391:THR:HG21	4:D:1501:HOH:O	2.03	0.59
1:A:177:LEU:HD13	1:A:183:LEU:HD11	1.84	0.59
1:F:54:TYR:N	4:F:1401:HOH:O	2.34	0.59
1:B:343:LEU:HA	1:B:377:ASN:HD21	1.67	0.59
1:A:414:PHE:CD2	1:A:427:LYS:HG2	2.38	0.59
1:B:343:LEU:HD13	1:B:376:GLU:HG3	1.83	0.59
1:A:177:LEU:HD13	1:A:183:LEU:CD1	2.33	0.59
1:D:17:SER:OG	1:D:90:HIS:HE1	1.85	0.59

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:ASN:ND2	1:B:319:ASP:H	1.98	0.58
1:E:343:LEU:HD11	1:E:376:GLU:HG3	1.84	0.58
1:D:343:LEU:CD1	1:D:376:GLU:HG3	2.33	0.58
1:B:259:LYS:HZ1	1:B:294:ASP:HB3	1.69	0.58
1:D:326:GLU:HG2	1:D:330:LYS:HE3	1.85	0.58
1:B:9:PHE:N	4:B:1412:HOH:O	2.35	0.58
1:B:60:TYR:O	1:B:158:PRO:HB2	2.03	0.58
1:E:33:VAL:H	1:E:136:ASN:HD21	1.50	0.58
1:D:88:LYS:HA	1:D:88:LYS:HE2	1.85	0.58
1:A:56:GLU:HG2	1:A:124:VAL:CG1	2.34	0.58
1:F:114:ILE:HG12	1:F:144:LEU:HD13	1.85	0.58
1:C:316:ASN:ND2	1:C:319:ASP:H	2.00	0.58
1:E:337:ASN:ND2	1:E:339:LYS:H	2.01	0.58
1:C:60:TYR:O	1:C:158:PRO:HB2	2.03	0.58
1:A:168:GLN:HG3	1:A:358:ILE:HD12	1.85	0.58
1:A:361:LEU:O	1:A:361:LEU:HD23	2.03	0.58
1:C:32:LYS:HD3	4:C:1394:HOH:O	2.03	0.58
1:A:168:GLN:HG3	1:A:358:ILE:CD1	2.33	0.58
1:F:234:LYS:HG3	1:F:474:ASP:OD2	2.04	0.57
1:D:270:SER:HA	1:D:302:ARG:NH2	2.19	0.57
1:C:259:LYS:HB2	1:C:259:LYS:HZ2	1.70	0.57
1:C:321:VAL:O	1:C:325:LEU:HG	2.05	0.57
1:F:158:PRO:HG3	1:F:206:ILE:CG2	2.35	0.57
1:D:155:SER:O	1:D:158:PRO:HD2	2.05	0.57
1:B:323:LYS:HA	1:B:326:GLU:HG2	1.87	0.57
1:B:392:ARG:HB3	1:B:392:ARG:NH1	2.20	0.57
1:A:155:SER:O	1:A:158:PRO:HD2	2.04	0.56
1:E:343:LEU:CD1	1:E:376:GLU:HG3	2.35	0.56
1:F:136:ASN:HD22	1:F:136:ASN:H	1.52	0.56
1:B:304:THR:HG22	1:B:346:PRO:HB2	1.87	0.56
1:B:418:VAL:HG13	1:B:419:ALA:N	2.20	0.56
1:A:182:ASN:C	1:A:183:LEU:HD12	2.25	0.56
1:D:265:ILE:CD1	1:E:25:GLN:HB3	2.36	0.56
1:F:305:GLN:HG3	4:F:1307:HOH:O	2.05	0.56
1:E:97:ASN:ND2	1:E:100:GLY:H	2.04	0.56
1:E:114:ILE:CG1	1:E:144:LEU:HD13	2.36	0.56
1:C:293:GLU:HG3	1:C:331:LYS:NZ	2.21	0.56
1:B:158:PRO:HG3	1:B:206:ILE:CG2	2.35	0.56
1:C:155:SER:O	1:C:158:PRO:HD2	2.06	0.56
1:F:299:ILE:HD12	1:F:308:LEU:HD22	1.87	0.56
1:B:478:LYS:HA	1:B:481:GLN:HE21	1.71	0.55

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:97:ASN:ND2	1:F:100:GLY:H	2.05	0.55
1:D:360:THR:O	1:D:364:ILE:HG13	2.06	0.55
1:D:285:ASN:OD1	1:D:289:LYS:HD2	2.07	0.55
1:A:60:TYR:O	1:A:158:PRO:HB2	2.05	0.55
1:F:236:PRO:HD2	4:F:994:HOH:O	2.05	0.55
1:E:149:GLU:HG3	1:E:399:PHE:CG	2.42	0.55
1:C:249:THR:CG2	1:C:265:ILE:HD11	2.36	0.55
1:B:17:SER:OG	1:B:90:HIS:HE1	1.90	0.55
1:F:364:ILE:O	1:F:368:MET:HG3	2.07	0.55
1:A:233:THR:HG22	1:A:473:PHE:HB3	1.89	0.55
1:F:17:SER:OG	1:F:90:HIS:HE1	1.90	0.55
1:C:62:LEU:HD22	1:C:66:LEU:HG	1.89	0.55
1:A:300:VAL:HG12	1:A:347:TYR:OH	2.06	0.55
1:A:25:GLN:HE21	1:C:249:THR:CG2	2.17	0.54
1:E:19:LYS:HA	1:E:22:HIS:CD2	2.42	0.54
1:F:431:SER:OG	1:F:433:HIS:HE1	1.90	0.54
1:E:33:VAL:H	1:E:136:ASN:ND2	2.05	0.54
1:A:90:HIS:HD2	1:C:240:TYR:HA	1.71	0.54
1:B:392:ARG:CB	1:B:392:ARG:HH11	2.20	0.54
1:F:344:LEU:N	1:F:377:ASN:HD21	1.88	0.54
1:C:259:LYS:HG2	4:C:1905:HOH:O	2.06	0.54
1:A:168:GLN:HG2	1:A:361:LEU:HD13	1.87	0.54
1:B:233:THR:HG22	1:B:473:PHE:HB3	1.88	0.54
1:D:334:VAL:HG22	4:D:1743:HOH:O	2.06	0.54
1:A:391:THR:CG2	1:A:393:ASP:H	2.18	0.54
1:A:56:GLU:HG2	4:A:1130:HOH:O	2.06	0.54
1:D:125:ILE:HD13	1:D:432:LEU:HD21	1.90	0.54
1:F:374:SER:OG	1:F:376:GLU:CG	2.53	0.54
1:A:57:THR:N	4:A:1130:HOH:O	2.41	0.54
1:C:149:GLU:OE2	1:C:399:PHE:HB3	2.07	0.54
1:C:359:ASN:O	1:C:363:GLU:HG3	2.07	0.54
1:B:112:LEU:O	1:B:464:ASN:HA	2.08	0.54
1:C:330:LYS:O	1:C:331:LYS:HD2	2.08	0.54
1:D:236:PRO:HD2	4:D:1061:HOH:O	2.06	0.54
1:C:56:GLU:HG2	1:C:124:VAL:CG1	2.38	0.54
1:B:391:THR:CG2	1:B:393:ASP:HB2	2.37	0.54
1:C:435:THR:HB	1:C:436:PRO:HD2	1.90	0.54
1:A:23:TYR:OH	1:A:24:LYS:HE2	2.08	0.54
1:B:472:SER:OG	1:B:475:GLU:HG3	2.08	0.54
1:E:174:LYS:O	1:E:178:GLU:HG3	2.08	0.54
1:D:399:PHE:O	1:D:400:LYS:HE3	2.08	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:391:THR:HG23	1:F:393:ASP:H	1.73	0.53
1:E:114:ILE:HD11	1:E:144:LEU:HB3	1.88	0.53
1:D:361:LEU:HD23	1:D:361:LEU:O	2.08	0.53
1:F:365:VAL:HA	1:F:368:MET:HE3	1.91	0.53
1:D:90:HIS:HD2	1:E:239:GLY:O	1.91	0.53
1:C:326:GLU:HG3	1:C:327:ILE:N	2.22	0.53
1:D:175:TYR:HB3	1:D:375:ILE:HG13	1.90	0.53
1:B:300:VAL:HG12	1:B:347:TYR:OH	2.08	0.53
1:C:286:ALA:N	4:C:1038:HOH:O	2.41	0.53
1:E:322:LEU:O	1:E:326:GLU:HG2	2.07	0.53
1:B:339:LYS:HA	1:B:339:LYS:HE2	1.90	0.53
1:C:374:SER:OG	1:C:376:GLU:HG2	2.08	0.53
1:A:283:ILE:HD13	1:A:283:ILE:O	2.08	0.53
1:F:41:GLU:HG3	1:F:128:GLY:HA3	1.91	0.53
1:D:32:LYS:HA	1:D:136:ASN:HD21	1.73	0.53
1:A:312:PRO:HG2	1:A:352:GLN:NE2	2.24	0.53
1:D:136:ASN:H	1:D:136:ASN:HD22	1.56	0.52
1:B:343:LEU:HD12	1:B:377:ASN:ND2	2.25	0.52
1:D:233:THR:CG2	1:D:473:PHE:HB3	2.39	0.52
1:D:265:ILE:HD12	1:E:25:GLN:HB3	1.91	0.52
1:A:168:GLN:NE2	1:A:190:LEU:HD11	2.24	0.52
1:D:21:THR:HG22	1:D:95:VAL:HG11	1.90	0.52
1:C:114:ILE:HD13	1:C:136:ASN:HA	1.91	0.52
1:B:21:THR:HG22	1:B:95:VAL:HG11	1.91	0.52
1:A:240:TYR:HA	1:C:90:HIS:HD2	1.73	0.52
1:E:29:ASN:N	1:E:29:ASN:ND2	2.58	0.52
1:D:300:VAL:HG12	1:D:347:TYR:OH	2.10	0.52
1:F:450:LEU:O	1:F:452:GLU:N	2.36	0.52
1:A:224:LEU:HD22	1:A:224:LEU:H	1.74	0.52
1:A:32:LYS:O	1:A:404:VAL:HG23	2.09	0.52
1:A:13:LEU:HD21	1:A:83:ALA:HA	1.92	0.52
1:A:155:SER:C	1:A:158:PRO:HD2	2.31	0.52
1:A:221:VAL:HA	1:A:224:LEU:HD23	1.92	0.51
1:A:27:PRO:HG3	1:C:253:TRP:CG	2.46	0.51
1:A:318:LEU:HD11	1:A:367:GLY:HA3	1.93	0.51
1:E:114:ILE:HD13	1:E:136:ASN:HA	1.93	0.51
1:D:62:LEU:HD22	1:D:66:LEU:HG	1.92	0.51
1:E:23:TYR:CE1	1:E:24:LYS:HG3	2.46	0.51
1:E:233:THR:HG22	1:E:473:PHE:HB3	1.91	0.51
1:A:418:VAL:HG13	1:A:419:ALA:N	2.26	0.51
1:F:466:LYS:HD3	4:F:848:HOH:O	2.10	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:446:GLY:C	1:C:448:GLY:H	2.12	0.51
1:A:317:PRO:O	1:A:321:VAL:HG23	2.11	0.51
1:F:434:ARG:HD2	4:F:1060:HOH:O	2.11	0.51
1:C:23:TYR:CE1	1:C:24:LYS:HG3	2.46	0.51
1:C:266:VAL:HG23	4:C:1561:HOH:O	2.09	0.51
1:B:392:ARG:HB3	1:B:392:ARG:HH11	1.76	0.51
1:F:400:LYS:CE	1:F:415:LYS:HD3	2.41	0.51
1:E:29:ASN:H	1:E:29:ASN:HD22	1.59	0.51
1:A:168:GLN:HG2	1:A:361:LEU:CD1	2.41	0.51
1:A:253:TRP:HZ2	1:A:264:HIS:HD1	1.59	0.51
1:C:418:VAL:HG13	1:C:419:ALA:N	2.26	0.51
1:B:63:GLN:HE22	1:B:471:TYR:N	1.94	0.50
1:F:123:PHE:HB3	1:F:125:ILE:CD1	2.40	0.50
3:D:504:NMN:H5R1	4:D:1482:HOH:O	2.11	0.50
1:D:113:PRO:HA	1:D:464:ASN:ND2	2.21	0.50
1:D:40:ARG:HD2	1:D:422:ASN:O	2.11	0.50
1:A:245:ALA:HB1	1:A:249:THR:CG2	2.42	0.50
1:A:172:LEU:HD13	1:A:189:LYS:HB2	1.92	0.50
1:F:151:ILE:HG23	1:F:152:LEU:N	2.27	0.50
1:D:344:LEU:N	1:D:377:ASN:HD21	2.03	0.50
1:C:155:SER:C	1:C:158:PRO:HD2	2.32	0.50
1:F:114:ILE:CG1	1:F:144:LEU:HD13	2.41	0.50
1:B:311:ARG:HB2	1:B:351:ILE:CG2	2.41	0.50
1:C:300:VAL:HG12	1:C:347:TYR:OH	2.11	0.50
1:E:333:PRO:O	1:E:345:PRO:HG3	2.12	0.50
1:E:257:HIS:HD2	4:E:781:HOH:O	1.94	0.50
1:A:249:THR:CG2	1:C:25:GLN:HE21	2.23	0.50
1:A:28:PRO:O	1:A:29:ASN:HB2	2.12	0.50
1:B:114:ILE:CG1	1:B:144:LEU:HD13	2.42	0.50
1:F:102:ASN:O	1:F:106:GLU:HG3	2.12	0.50
1:F:400:LYS:HE2	1:F:415:LYS:HD3	1.94	0.50
1:A:404:VAL:C	4:A:1330:HOH:O	2.50	0.50
1:B:97:ASN:ND2	1:B:100:GLY:H	2.10	0.50
1:C:258:GLU:OE2	1:C:281:TYR:HD1	1.94	0.50
1:C:202:GLU:O	1:C:206:ILE:HG13	2.11	0.50
1:B:361:LEU:O	1:B:365:VAL:HG22	2.12	0.50
1:D:224:LEU:H	1:D:224:LEU:HD22	1.77	0.50
1:F:68:LYS:HA	4:F:2060:HOH:O	2.12	0.49
1:C:11:ILE:HG23	1:C:12:LEU:HD12	1.94	0.49
1:E:335:THR:OG1	1:E:336:GLU:N	2.45	0.49
1:C:120:PRO:O	1:C:123:PHE:HB2	2.12	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:TYR:CE2	1:E:97:ASN:HB2	2.47	0.49
1:F:37:PHE:CZ	1:F:397:CYS:HB3	2.47	0.49
1:D:257:HIS:HD2	1:D:260:ASP:OD2	1.96	0.49
1:A:253:TRP:CD1	1:C:27:PRO:HG3	2.47	0.49
1:A:444:GLU:O	1:A:447:LYS:HG3	2.13	0.49
1:C:33:VAL:H	1:C:136:ASN:ND2	2.11	0.49
1:A:264:HIS:NE2	1:A:268:GLN:NE2	2.60	0.49
1:E:136:ASN:ND2	1:E:136:ASN:H	2.10	0.49
1:F:63:GLN:NE2	1:F:471:TYR:H	1.97	0.49
1:E:60:TYR:O	1:E:158:PRO:HB2	2.13	0.49
1:B:323:LYS:O	1:B:327:ILE:HG12	2.12	0.49
1:E:360:THR:O	1:E:364:ILE:HG13	2.13	0.49
1:F:19:LYS:HA	1:F:22:HIS:CD2	2.48	0.49
1:D:63:GLN:HE22	1:D:471:TYR:N	2.02	0.49
1:F:32:LYS:HA	1:F:136:ASN:HD21	1.77	0.49
1:D:253:TRP:CD2	1:E:27:PRO:HG3	2.48	0.49
1:D:339:LYS:NZ	1:D:339:LYS:HA	2.28	0.49
1:A:318:LEU:HG	1:A:364:ILE:HA	1.95	0.48
1:A:19:LYS:HA	1:A:22:HIS:CD2	2.47	0.48
1:F:40:ARG:HD2	1:F:422:ASN:O	2.13	0.48
1:A:432:LEU:HD11	1:A:440:PHE:HB3	1.96	0.48
1:C:114:ILE:CG1	1:C:144:LEU:HD13	2.43	0.48
1:D:90:HIS:CD2	1:E:240:TYR:HA	2.47	0.48
1:E:290:ILE:HD11	4:E:1426:HOH:O	2.12	0.48
1:C:34:TYR:HB3	1:C:403:TYR:HB3	1.95	0.48
1:A:97:ASN:ND2	1:A:100:GLY:H	2.12	0.48
1:B:34:TYR:HB3	1:B:403:TYR:HB3	1.95	0.48
1:A:413:VAL:HG11	1:C:252:ALA:HA	1.94	0.48
1:F:288:GLU:HG3	1:F:331:LYS:HE3	1.96	0.48
1:E:183:LEU:HD11	1:E:484:ILE:HG21	1.94	0.48
1:A:183:LEU:HD21	1:A:484:ILE:HG21	1.95	0.48
1:E:63:GLN:NE2	1:E:471:TYR:H	2.04	0.48
1:C:236:PRO:HD2	4:C:1302:HOH:O	2.13	0.48
1:C:472:SER:OG	1:C:475:GLU:HG3	2.14	0.48
1:F:21:THR:HG22	1:F:95:VAL:HG21	1.94	0.48
1:B:155:SER:O	1:B:158:PRO:HD2	2.14	0.48
1:E:34:TYR:HB3	1:E:403:TYR:HB3	1.95	0.48
1:D:427:LYS:HE3	4:D:1275:HOH:O	2.13	0.48
1:E:17:SER:OG	1:E:90:HIS:HE1	1.96	0.48
1:C:452:GLU:HG3	4:E:1169:HOH:O	2.13	0.48
1:A:114:ILE:O	1:A:463:LYS:HA	2.13	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:TYR:O	1:A:127:ARG:CD	2.61	0.47
1:A:17:SER:OG	1:A:90:HIS:HE1	1.97	0.47
1:D:55:GLU:HB2	4:D:1190:HOH:O	2.13	0.47
1:A:62:LEU:HD22	1:A:66:LEU:HG	1.95	0.47
1:B:228:LYS:HE3	1:B:228:LYS:HB2	1.66	0.47
1:C:277:VAL:HG13	1:C:277:VAL:O	2.14	0.47
1:C:311:ARG:HH21	3:C:503:NMN:H2RC	1.78	0.47
1:C:127:ARG:NH2	1:C:396:ASN:HD22	2.11	0.47
1:D:450:LEU:HB2	1:D:452:GLU:HG2	1.95	0.47
1:C:224:LEU:CD1	1:C:238:PRO:HD2	2.42	0.47
1:C:114:ILE:O	1:C:463:LYS:HA	2.15	0.47
1:F:136:ASN:ND2	1:F:136:ASN:H	2.12	0.47
1:C:55:GLU:O	1:C:56:GLU:HB2	2.13	0.47
1:D:80:ILE:HD13	1:D:102:ASN:ND2	2.29	0.47
1:E:311:ARG:NH1	3:E:505:NMN:O2R	2.45	0.47
1:D:56:GLU:CG	1:D:124:VAL:HG12	2.43	0.47
1:B:351:ILE:HG23	1:B:351:ILE:O	2.13	0.47
1:D:339:LYS:HZ1	1:D:339:LYS:HA	1.79	0.47
1:D:77:LYS:HD2	1:D:77:LYS:HA	1.72	0.47
1:D:60:TYR:O	1:D:158:PRO:HB2	2.15	0.47
1:B:478:LYS:HG3	1:B:481:GLN:NE2	2.28	0.47
1:A:312:PRO:HG2	1:A:352:GLN:HE22	1.79	0.47
1:C:224:LEU:H	1:C:224:LEU:HD22	1.79	0.47
1:A:32:LYS:HB3	1:A:405:VAL:HB	1.97	0.47
1:C:446:GLY:C	1:C:448:GLY:N	2.68	0.47
1:C:266:VAL:O	1:C:302:ARG:NH2	2.47	0.47
1:A:339:LYS:HB3	4:A:1129:HOH:O	2.13	0.47
1:B:250:ILE:HD13	1:B:262:PHE:CE1	2.44	0.47
1:B:120:PRO:HB2	1:B:123:PHE:CD1	2.50	0.47
1:A:344:LEU:HG	1:A:377:ASN:ND2	2.30	0.47
1:F:391:THR:HG22	1:F:393:ASP:H	1.77	0.47
1:B:392:ARG:NE	4:B:812:HOH:O	2.36	0.47
1:D:221:VAL:O	1:D:224:LEU:HD23	2.14	0.47
1:B:19:LYS:HA	1:B:22:HIS:CD2	2.49	0.47
1:E:120:PRO:O	1:E:123:PHE:HB2	2.15	0.47
1:F:226:LEU:HD23	1:F:226:LEU:C	2.35	0.47
1:A:158:PRO:HG3	1:A:206:ILE:CG2	2.45	0.46
1:D:158:PRO:HG3	1:D:206:ILE:CG2	2.45	0.46
1:F:447:LYS:HE2	4:F:1479:HOH:O	2.15	0.46
1:A:291:TRP:HE3	1:A:299:ILE:HD11	1.77	0.46
1:B:418:VAL:HG13	1:B:419:ALA:H	1.79	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:LEU:HD22	1:B:66:LEU:HG	1.97	0.46
1:E:119:VAL:HG21	1:E:125:ILE:CD1	2.32	0.46
1:B:236:PRO:HD2	4:B:863:HOH:O	2.13	0.46
1:E:431:SER:OG	1:E:433:HIS:HE1	1.98	0.46
1:E:316:ASN:HD22	1:E:319:ASP:N	2.02	0.46
1:A:146:ASN:ND2	1:C:246:GLU:OE2	2.40	0.46
1:F:414:PHE:CD2	1:F:427:LYS:HE2	2.51	0.46
1:A:31:SER:HB2	1:A:139:PRO:HB3	1.96	0.46
1:A:320:THR:O	1:A:324:VAL:HG23	2.16	0.46
1:C:172:LEU:HD13	1:C:189:LYS:HB2	1.97	0.46
1:F:444:GLU:O	1:F:447:LYS:HG3	2.16	0.46
1:F:443:LEU:HD12	1:F:448:GLY:HA2	1.98	0.46
1:D:97:ASN:ND2	1:D:100:GLY:H	2.12	0.46
1:F:418:VAL:HG13	1:F:419:ALA:N	2.29	0.46
1:C:439:ASN:OD1	1:C:440:PHE:N	2.43	0.46
1:C:12:LEU:HD23	1:C:96:PHE:HZ	1.81	0.46
1:B:32:LYS:HB3	1:B:405:VAL:HB	1.97	0.46
1:A:119:VAL:HG21	1:A:125:ILE:CD1	2.44	0.46
1:C:172:LEU:HD13	1:C:189:LYS:CB	2.46	0.46
1:A:351:ILE:HG23	1:A:351:ILE:O	2.15	0.46
1:A:188:TYR:HA	4:A:1001:HOH:O	2.16	0.46
1:F:450:LEU:HB2	1:F:452:GLU:HG3	1.96	0.46
1:B:76:THR:O	1:B:80:ILE:HG13	2.16	0.46
1:A:334:VAL:HG22	1:A:344:LEU:HD23	1.97	0.46
1:A:56:GLU:HG2	1:A:124:VAL:HG12	1.98	0.46
1:D:302:ARG:HD3	1:D:306:ALA:O	2.15	0.46
1:E:262:PHE:CD2	1:E:290:ILE:HG21	2.50	0.46
1:F:375:ILE:O	1:F:375:ILE:HG22	2.16	0.46
1:C:9:PHE:N	4:C:2039:HOH:O	2.48	0.46
1:E:337:ASN:HD22	1:E:339:LYS:H	1.64	0.45
1:D:239:GLY:O	1:E:90:HIS:HD2	1.99	0.45
1:C:366:GLU:HG3	1:C:370:GLN:HE21	1.81	0.45
1:C:431:SER:OG	1:C:433:HIS:HE1	1.98	0.45
1:B:291:TRP:HE3	1:B:299:ILE:HD11	1.82	0.45
1:A:119:VAL:CG2	1:A:120:PRO:HD2	2.46	0.45
1:C:318:LEU:HD23	1:C:364:ILE:HA	1.97	0.45
1:F:155:SER:O	1:F:158:PRO:HD2	2.17	0.45
1:A:233:THR:CG2	1:A:473:PHE:HB3	2.46	0.45
1:D:9:PHE:N	4:D:1096:HOH:O	2.50	0.45
1:C:194:GLY:N	1:C:383:GLY:HA2	2.31	0.45
1:E:13:LEU:HD11	1:E:82:GLU:OE2	2.17	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:LYS:HB3	4:D:2028:HOH:O	2.16	0.45
1:A:333:PRO:O	1:A:345:PRO:HG3	2.16	0.45
1:A:149:GLU:OE2	1:A:399:PHE:HB3	2.15	0.45
1:C:226:LEU:HD23	1:C:226:LEU:C	2.36	0.45
1:A:119:VAL:HG23	1:A:120:PRO:HD2	1.98	0.45
1:C:247:HIS:CE1	1:C:277:VAL:HG21	2.51	0.45
1:F:54:TYR:CE1	1:F:394:LEU:HD11	2.52	0.45
1:C:293:GLU:HG3	1:C:331:LYS:HZ3	1.81	0.45
1:C:56:GLU:HG2	1:C:124:VAL:HG12	1.99	0.45
1:B:300:VAL:HG12	1:B:347:TYR:CZ	2.52	0.45
1:F:33:VAL:H	1:F:136:ASN:HD21	1.65	0.45
1:A:253:TRP:CG	1:C:27:PRO:HG3	2.52	0.45
1:D:467:VAL:HG11	4:D:1532:HOH:O	2.16	0.45
1:A:221:VAL:HG12	4:A:1786:HOH:O	2.17	0.45
1:D:112:LEU:O	1:D:464:ASN:HA	2.17	0.45
1:F:484:ILE:HG13	4:F:1506:HOH:O	2.17	0.45
1:C:33:VAL:H	1:C:136:ASN:HD21	1.64	0.45
1:A:183:LEU:HD12	1:A:183:LEU:N	2.31	0.45
1:F:114:ILE:HD13	1:F:136:ASN:HA	1.97	0.45
1:A:23:TYR:CE2	1:A:97:ASN:HB2	2.52	0.45
1:C:284:TYR:C	1:C:286:ALA:H	2.20	0.45
1:A:288:GLU:OE2	1:E:436:PRO:HD2	2.17	0.45
1:A:179:THR:HG22	1:A:341:TYR:CE2	2.52	0.45
1:C:171:ILE:HG23	1:C:362:GLN:NE2	2.32	0.45
1:F:316:ASN:HD22	1:F:319:ASP:N	2.04	0.45
1:C:343:LEU:HA	1:C:377:ASN:HD21	1.82	0.45
1:D:300:VAL:HA	1:D:347:TYR:CZ	2.51	0.45
1:C:285:ASN:C	4:C:1038:HOH:O	2.56	0.45
1:C:174:LYS:O	1:C:178:GLU:HG3	2.17	0.45
1:D:431:SER:OG	1:D:433:HIS:HE1	2.00	0.45
1:A:226:LEU:C	1:A:226:LEU:HD23	2.38	0.45
1:D:24:LYS:O	1:E:264:HIS:HE1	1.98	0.45
1:B:112:LEU:HA	1:B:113:PRO:HD3	1.82	0.44
1:E:155:SER:C	1:E:158:PRO:HD2	2.36	0.44
1:C:168:GLN:HG2	1:C:358:ILE:HD12	1.99	0.44
1:D:375:ILE:HG22	1:D:375:ILE:O	2.17	0.44
1:A:262:PHE:O	1:A:266:VAL:HG13	2.17	0.44
1:E:228:LYS:HB2	4:E:2059:HOH:O	2.17	0.44
1:D:144:LEU:HD22	4:D:1329:HOH:O	2.16	0.44
1:D:302:ARG:HB2	1:D:347:TYR:HB2	1.99	0.44
1:C:445:GLU:HA	4:C:849:HOH:O	2.17	0.44

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:ILE:HG21	1:A:380:PHE:HE2	1.82	0.44
1:B:237:VAL:HA	1:B:238:PRO:HD3	1.81	0.44
1:E:414:PHE:CE2	1:E:427:LYS:HE3	2.53	0.44
1:E:324:VAL:O	1:E:328:LEU:HG	2.18	0.44
1:B:226:LEU:HD23	1:B:226:LEU:C	2.38	0.44
1:C:224:LEU:N	1:C:224:LEU:HD22	2.32	0.44
1:D:224:LEU:HD12	1:D:238:PRO:HD2	1.98	0.44
1:B:191:HIS:CD2	1:B:193:PHE:HE2	2.36	0.44
1:A:311:ARG:HH22	2:A:601:PO4:P	2.40	0.44
1:C:291:TRP:HE3	1:C:299:ILE:HD11	1.79	0.44
1:B:323:LYS:O	1:B:326:GLU:HG2	2.18	0.44
1:F:378:ILE:CG1	1:F:379:ALA:N	2.81	0.44
1:B:324:VAL:O	1:B:328:LEU:HG	2.18	0.44
1:C:404:VAL:HG12	1:C:405:VAL:N	2.33	0.44
1:A:82:GLU:O	1:A:86:VAL:HG23	2.17	0.44
1:C:437:ALA:HB3	1:E:336:GLU:OE2	2.17	0.44
1:C:24:LYS:HZ3	1:C:95:VAL:HG13	1.83	0.44
1:A:410:GLY:HA2	4:A:1330:HOH:O	2.16	0.44
1:D:32:LYS:HB3	1:D:405:VAL:HB	1.99	0.44
1:C:172:LEU:HD21	1:C:361:LEU:HD21	1.98	0.44
1:E:375:ILE:HG22	1:E:375:ILE:O	2.17	0.44
1:A:376:GLU:HG2	1:A:376:GLU:H	1.62	0.44
1:A:478:LYS:HA	1:A:481:GLN:HE21	1.82	0.44
1:B:114:ILE:HD11	1:B:144:LEU:HB3	2.01	0.43
1:B:233:THR:CG2	1:B:473:PHE:HB3	2.48	0.43
1:D:229:LYS:HD3	1:D:230:TYR:CE1	2.53	0.43
1:C:177:LEU:HD13	1:C:183:LEU:CD1	2.48	0.43
1:B:433:HIS:HD2	1:B:454:GLY:O	2.00	0.43
1:C:420:ASP:OD1	1:C:422:ASN:HB2	2.18	0.43
1:B:378:ILE:HG12	1:B:379:ALA:N	2.33	0.43
1:F:125:ILE:N	1:F:125:ILE:HD12	2.33	0.43
1:D:136:ASN:H	1:D:136:ASN:ND2	2.17	0.43
1:D:326:GLU:O	1:D:330:LYS:HG3	2.17	0.43
1:C:443:LEU:HD12	1:C:448:GLY:HA2	2.01	0.43
1:B:351:ILE:HD13	1:B:351:ILE:C	2.37	0.43
1:D:238:PRO:HA	4:D:1452:HOH:O	2.17	0.43
1:F:183:LEU:HD11	1:F:484:ILE:CG2	2.46	0.43
1:A:361:LEU:C	1:A:361:LEU:HD23	2.39	0.43
1:A:283:ILE:HD13	1:A:287:CYS:SG	2.59	0.43
1:F:40:ARG:NH2	4:F:841:HOH:O	2.49	0.43
1:F:9:PHE:N	4:F:1116:HOH:O	2.52	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:ARG:NH2	1:C:396:ASN:ND2	2.66	0.43
1:B:333:PRO:O	1:B:345:PRO:HG3	2.18	0.43
1:C:375:ILE:C	1:C:377:ASN:H	2.22	0.43
1:D:282:ASP:HB3	1:D:285:ASN:HB3	2.01	0.43
1:D:105:LEU:O	1:D:109:ASP:HA	2.19	0.43
1:F:274:VAL:O	1:F:274:VAL:HG13	2.18	0.43
1:D:351:ILE:O	1:D:351:ILE:HG23	2.19	0.43
1:A:344:LEU:HG	1:A:377:ASN:HD22	1.82	0.43
1:D:63:GLN:NE2	1:D:471:TYR:H	2.02	0.43
1:F:60:TYR:O	1:F:158:PRO:HB2	2.18	0.43
1:E:149:GLU:HG3	1:E:399:PHE:CD2	2.53	0.43
1:A:247:HIS:CE1	1:A:277:VAL:HG21	2.54	0.43
1:E:296:ARG:HG2	4:E:1948:HOH:O	2.18	0.43
1:C:264:HIS:CE1	1:C:268:GLN:HG3	2.54	0.43
1:D:391:THR:CG2	1:D:392:ARG:N	2.81	0.43
1:C:246:GLU:O	1:C:249:THR:HB	2.19	0.43
1:F:391:THR:HG21	4:F:1253:HOH:O	2.17	0.43
1:C:191:HIS:CD2	1:C:193:PHE:HE2	2.37	0.43
1:B:182:ASN:HB2	1:B:184:ASP:OD1	2.19	0.43
1:B:391:THR:HG21	1:B:393:ASP:HB2	2.00	0.43
1:C:311:ARG:HH22	2:C:605:PO4:P	2.42	0.43
1:C:360:THR:O	1:C:364:ILE:HG13	2.19	0.43
1:A:304:THR:HG23	1:A:346:PRO:HB2	2.00	0.43
1:D:19:LYS:HA	1:D:22:HIS:CD2	2.54	0.43
1:C:322:LEU:CD2	1:C:371:LYS:HG3	2.44	0.42
1:E:337:ASN:C	1:E:337:ASN:ND2	2.70	0.42
1:C:288:GLU:N	4:C:1038:HOH:O	2.51	0.42
1:A:114:ILE:CG2	1:A:115:GLU:N	2.82	0.42
1:B:188:TYR:HA	4:B:1681:HOH:O	2.19	0.42
1:B:183:LEU:HD11	1:B:484:ILE:HD12	2.01	0.42
1:A:134:VAL:HG11	1:A:148:ILE:HD11	2.01	0.42
1:F:33:VAL:H	1:F:136:ASN:ND2	2.17	0.42
1:B:113:PRO:HD2	1:B:144:LEU:HD12	2.01	0.42
1:D:198:VAL:HG13	1:D:200:SER:H	1.84	0.42
1:D:73:LYS:HE2	1:D:76:THR:HG22	2.02	0.42
1:B:257:HIS:HD2	1:B:260:ASP:OD1	2.03	0.42
1:A:73:LYS:HE3	1:A:76:THR:HG22	2.00	0.42
1:B:77:LYS:CD	1:B:105:LEU:HD21	2.43	0.42
1:C:158:PRO:HG3	1:C:206:ILE:CG2	2.49	0.42
1:B:392:ARG:NH2	4:B:812:HOH:O	2.46	0.42
1:A:172:LEU:HD13	1:A:189:LYS:CB	2.48	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:27:PRO:HG3	1:E:253:TRP:CG	2.54	0.42
1:E:337:ASN:HB2	4:E:1717:HOH:O	2.19	0.42
1:A:451:GLU:N	4:A:1898:HOH:O	2.53	0.42
1:F:302:ARG:HB2	1:F:347:TYR:HB2	2.01	0.42
1:C:351:ILE:HG23	1:C:351:ILE:O	2.19	0.42
1:B:326:GLU:HG3	1:B:327:ILE:N	2.34	0.42
1:C:358:ILE:HG23	1:C:359:ASN:N	2.35	0.42
1:A:283:ILE:HD13	1:A:287:CYS:HG	1.84	0.42
1:F:226:LEU:HD23	1:F:226:LEU:O	2.20	0.42
1:C:409:LEU:HD13	1:C:410:GLY:N	2.35	0.42
1:C:78:GLU:H	1:C:78:GLU:CD	2.23	0.42
1:F:368:MET:CE	1:F:378:ILE:HG21	2.49	0.42
1:F:116:ILE:HB	1:F:462:PHE:HB3	2.01	0.42
1:A:406:THR:O	1:A:407:ASN:HB2	2.19	0.42
1:A:245:ALA:HB1	1:A:249:THR:HG22	2.02	0.42
1:B:430:LEU:HD22	1:B:444:GLU:HG2	2.00	0.42
1:B:305:GLN:HE21	1:B:305:GLN:HB3	1.48	0.42
1:B:351:ILE:HD13	1:B:352:GLN:N	2.35	0.42
1:D:224:LEU:N	1:D:224:LEU:HD22	2.35	0.42
1:C:378:ILE:HG12	1:C:379:ALA:N	2.35	0.42
1:B:444:GLU:O	1:B:447:LYS:HG3	2.20	0.42
1:C:450:LEU:O	1:C:452:GLU:N	2.48	0.42
1:B:191:HIS:CD2	1:B:193:PHE:CE2	3.08	0.42
1:C:138:ASP:HA	1:C:139:PRO:HD2	1.87	0.42
1:F:333:PRO:O	1:F:345:PRO:HG3	2.19	0.42
1:E:170:LYS:HG2	1:E:482:LEU:HD22	2.02	0.42
1:A:284:TYR:CZ	1:A:323:LYS:HD3	2.55	0.42
1:F:155:SER:C	1:F:158:PRO:HD2	2.40	0.41
1:A:283:ILE:HD12	1:A:324:VAL:CG2	2.50	0.41
1:F:351:ILE:O	1:F:351:ILE:HG23	2.20	0.41
1:C:318:LEU:C	1:C:318:LEU:HD13	2.40	0.41
1:C:434:ARG:HD2	1:C:438:GLY:HA2	2.02	0.41
1:A:202:GLU:HG2	1:C:200:SER:HB2	2.02	0.41
1:D:316:ASN:HD22	1:D:319:ASP:N	1.98	0.41
1:B:155:SER:C	1:B:158:PRO:HD2	2.40	0.41
1:A:176:LEU:HD23	1:A:183:LEU:HA	2.03	0.41
1:B:21:THR:HG22	1:B:95:VAL:CG1	2.49	0.41
1:C:375:ILE:C	1:C:377:ASN:N	2.74	0.41
1:E:11:ILE:HG23	1:E:12:LEU:HD12	2.01	0.41
1:E:469:LYS:HE3	1:E:471:TYR:CE2	2.55	0.41
1:B:291:TRP:CE3	1:B:299:ILE:HD11	2.55	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:GLY:HA2	4:C:977:HOH:O	2.20	0.41
1:D:197:GLY:HA2	1:E:392:ARG:CD	2.50	0.41
1:D:311:ARG:HH22	2:D:607:PO4:P	2.43	0.41
1:E:423:LYS:HE3	4:E:1606:HOH:O	2.21	0.41
1:C:348:LEU:O	1:C:349:ARG:NH1	2.46	0.41
1:E:137:THR:OG1	1:E:464:ASN:ND2	2.50	0.41
1:B:285:ASN:O	1:B:289:LYS:HB3	2.19	0.41
1:B:259:LYS:CB	1:B:259:LYS:NZ	2.80	0.41
1:D:155:SER:C	1:D:158:PRO:HD2	2.41	0.41
1:C:288:GLU:OE2	1:F:436:PRO:HD2	2.21	0.41
1:F:105:LEU:O	1:F:109:ASP:HA	2.20	0.41
1:A:316:ASN:HB3	1:A:319:ASP:HB2	2.02	0.41
1:A:224:LEU:HD22	1:A:224:LEU:N	2.35	0.41
1:A:391:THR:HG23	1:A:392:ARG:N	2.35	0.41
1:B:427:LYS:O	1:B:444:GLU:HB3	2.20	0.41
1:F:303:SER:OG	1:F:305:GLN:NE2	2.54	0.41
1:C:266:VAL:HB	1:C:302:ARG:HH21	1.85	0.41
1:D:312:PRO:HD2	1:D:351:ILE:O	2.21	0.41
1:E:259:LYS:HD2	1:E:294:ASP:HB3	2.02	0.41
1:B:178:GLU:OE1	1:B:369:LYS:HE3	2.21	0.41
1:D:262:PHE:O	1:D:266:VAL:HG22	2.20	0.41
1:A:127:ARG:NH2	1:A:396:ASN:HD22	2.19	0.41
1:B:156:TRP:CG	1:B:392:ARG:HG3	2.56	0.41
1:F:28:PRO:HG3	4:F:1730:HOH:O	2.21	0.41
1:A:237:VAL:HA	1:A:238:PRO:HD3	1.79	0.40
1:A:54:TYR:N	4:A:1619:HOH:O	2.54	0.40
1:F:114:ILE:O	1:F:463:LYS:HA	2.21	0.40
1:C:284:TYR:CE2	1:C:323:LYS:HD3	2.56	0.40
1:C:443:LEU:HD22	1:E:372:MET:HE3	2.01	0.40
1:C:362:GLN:O	1:C:365:VAL:HG22	2.21	0.40
1:C:404:VAL:O	1:C:410:GLY:HA2	2.21	0.40
1:C:273:PRO:HG3	1:C:306:ALA:HA	2.03	0.40
1:B:262:PHE:CD2	1:B:291:TRP:CZ2	3.09	0.40
1:C:120:PRO:HB2	1:C:123:PHE:CD1	2.56	0.40
1:C:324:VAL:O	1:C:328:LEU:HG	2.21	0.40
1:A:75:VAL:HB	1:A:110:GLY:HA2	2.03	0.40
1:A:249:THR:HG21	1:C:25:GLN:HE22	1.79	0.40
1:C:259:LYS:NZ	1:C:259:LYS:CB	2.84	0.40
1:B:306:ALA:N	1:B:307:PRO:CD	2.82	0.40
1:D:236:PRO:HB3	1:E:89:GLU:HG3	2.04	0.40
1:C:198:VAL:CG1	1:C:199:SER:N	2.84	0.40

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LYS:O	1:A:293:GLU:HB2	2.22	0.40
1:C:437:ALA:HB3	1:E:336:GLU:CD	2.41	0.40
1:F:124:VAL:C	1:F:125:ILE:HD12	2.41	0.40
1:B:346:PRO:N	4:B:2001:HOH:O	2.53	0.40
1:A:62:LEU:O	1:A:66:LEU:HG	2.22	0.40
1:B:366:GLU:HG3	1:B:370:GLN:HE21	1.86	0.40
1:A:219:ASP:CG	1:C:17:SER:HB2	2.42	0.40
1:A:350:VAL:HG12	1:A:351:ILE:N	2.36	0.40
1:E:370:GLN:CD	4:E:1418:HOH:O	2.59	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

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

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	455/491 (93%)	424 (93%)	30 (7%)	1 (0%)	52	59
1	B	460/491 (94%)	430 (94%)	29 (6%)	1 (0%)	52	59
1	C	455/491 (93%)	425 (93%)	26 (6%)	4 (1%)	21	19
1	D	460/491 (94%)	444 (96%)	15 (3%)	1 (0%)	52	59
1	E	460/491 (94%)	442 (96%)	16 (4%)	2 (0%)	39	42
1	F	457/491 (93%)	438 (96%)	18 (4%)	1 (0%)	52	59
All	All	2747/2946 (93%)	2603 (95%)	134 (5%)	10 (0%)	39	42

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	483	ASN
1	D	454	GLY
1	F	451	GLU

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	56	GLU
1	E	451	GLU
1	B	97	ASN
1	C	451	GLU
1	E	454	GLY
1	A	451	GLU
1	C	40	ARG

### 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	405/430 (94%)	389 (96%)	16 (4%)	38	47
1	B	408/430 (95%)	391 (96%)	17 (4%)	36	44
1	C	405/430 (94%)	393 (97%)	12 (3%)	48	60
1	D	408/430 (95%)	388 (95%)	20 (5%)	31	36
1	E	408/430 (95%)	390 (96%)	18 (4%)	35	42
1	F	407/430 (95%)	393 (97%)	14 (3%)	44	54
All	All	2441/2580 (95%)	2344 (96%)	97 (4%)	38	47

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	TYR
1	A	62	LEU
1	A	94	ASP
1	A	97	ASN
1	A	144	LEU
1	A	190	LEU
1	A	195	TYR
1	A	219	ASP
1	A	249	THR
1	A	283	ILE
1	A	294	ASP

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	318	LEU
1	A	335	THR
1	A	370	GLN
1	A	391	THR
1	A	398	SER
1	B	18	TYR
1	B	32	LYS
1	B	62	LEU
1	B	77	LYS
1	B	97	ASN
1	B	144	LEU
1	B	190	LEU
1	B	195	TYR
1	B	219	ASP
1	B	249	THR
1	B	259	LYS
1	B	294	ASP
1	B	305	GLN
1	B	318	LEU
1	B	351	ILE
1	B	391	THR
1	B	483	ASN
1	C	18	TYR
1	C	62	LEU
1	C	97	ASN
1	C	123	PHE
1	C	144	LEU
1	C	195	TYR
1	C	198	VAL
1	C	219	ASP
1	C	249	THR
1	C	259	LYS
1	C	294	ASP
1	C	386	LEU
1	D	18	TYR
1	D	62	LEU
1	D	88	LYS
1	D	97	ASN
1	D	136	ASN
1	D	144	LEU
1	D	183	LEU
1	D	186	LEU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	190	LEU
1	D	195	TYR
1	D	198	VAL
1	D	219	ASP
1	D	249	THR
1	D	318	LEU
1	D	339	LYS
1	D	372	MET
1	D	400	LYS
1	D	434	ARG
1	D	455	GLN
1	D	467	VAL
1	E	18	TYR
1	E	29	ASN
1	E	62	LEU
1	E	97	ASN
1	E	102	ASN
1	E	136	ASN
1	E	144	LEU
1	E	186	LEU
1	E	195	TYR
1	E	219	ASP
1	E	249	THR
1	E	275	SER
1	E	289	LYS
1	E	310	ILE
1	E	318	LEU
1	E	337	ASN
1	E	386	LEU
1	E	409	LEU
1	F	18	TYR
1	F	62	LEU
1	F	97	ASN
1	F	136	ASN
1	F	144	LEU
1	F	190	LEU
1	F	195	TYR
1	F	219	ASP
1	F	290	ILE
1	F	318	LEU
1	F	391	THR
1	F	394	LEU

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	F	404	VAL
1	F	409	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (121) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	63	GLN
1	A	90	HIS
1	A	97	ASN
1	A	102	ASN
1	A	129	ASN
1	A	136	ASN
1	A	168	GLN
1	A	201	GLN
1	A	268	GLN
1	A	316	ASN
1	A	352	GLN
1	A	362	GLN
1	A	370	GLN
1	A	377	ASN
1	A	396	ASN
1	A	433	HIS
1	A	464	ASN
1	A	479	ASN
1	A	481	GLN
1	B	25	GLN
1	B	63	GLN
1	B	90	HIS
1	B	92	GLN
1	B	97	ASN
1	B	102	ASN
1	B	136	ASN
1	B	201	GLN
1	B	257	HIS
1	B	305	GLN
1	B	316	ASN
1	B	337	ASN
1	B	362	GLN
1	B	370	GLN
1	B	377	ASN
1	B	396	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	433	HIS
1	B	459	HIS
1	B	464	ASN
1	B	479	ASN
1	B	481	GLN
1	C	22	HIS
1	C	25	GLN
1	C	29	ASN
1	C	63	GLN
1	C	81	GLN
1	C	90	HIS
1	C	97	ASN
1	C	136	ASN
1	C	201	GLN
1	C	264	HIS
1	C	285	ASN
1	C	316	ASN
1	C	362	GLN
1	C	370	GLN
1	C	377	ASN
1	C	396	ASN
1	C	407	ASN
1	C	433	HIS
1	C	464	ASN
1	C	479	ASN
1	C	481	GLN
1	C	483	ASN
1	D	25	GLN
1	D	29	ASN
1	D	63	GLN
1	D	90	HIS
1	D	97	ASN
1	D	102	ASN
1	D	129	ASN
1	D	136	ASN
1	D	201	GLN
1	D	214	ASN
1	D	257	HIS
1	D	264	HIS
1	D	268	GLN
1	D	316	ASN
1	D	362	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	377	ASN
1	D	396	ASN
1	D	433	HIS
1	D	464	ASN
1	D	479	ASN
1	E	22	HIS
1	E	25	GLN
1	E	29	ASN
1	E	63	GLN
1	E	90	HIS
1	E	97	ASN
1	E	102	ASN
1	E	136	ASN
1	E	201	GLN
1	E	257	HIS
1	E	264	HIS
1	E	316	ASN
1	E	337	ASN
1	E	352	GLN
1	E	362	GLN
1	E	370	GLN
1	E	377	ASN
1	E	433	HIS
1	E	464	ASN
1	E	479	ASN
1	F	22	HIS
1	F	25	GLN
1	F	63	GLN
1	F	90	HIS
1	F	97	ASN
1	F	136	ASN
1	F	201	GLN
1	F	257	HIS
1	F	305	GLN
1	F	316	ASN
1	F	362	GLN
1	F	377	ASN
1	F	396	ASN
1	F	433	HIS
1	F	439	ASN
1	F	464	ASN
1	F	479	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	F	481	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 ⓘ

18 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	$\# Z  > 2$
3	NMN	A	501	-	20,23,23	3.38	10 (50%)	26,34,34	3.51	10 (38%)
2	PO4	A	601	-	4,4,4	1.01	0	6,6,6	0.27	0
2	PO4	A	606	-	4,4,4	1.12	0	6,6,6	0.27	0
3	NMN	B	502	-	20,23,23	3.38	10 (50%)	26,34,34	3.41	10 (38%)
2	PO4	B	603	-	4,4,4	1.14	0	6,6,6	0.27	0
2	PO4	B	604	-	4,4,4	1.13	0	6,6,6	0.27	0
3	NMN	C	503	-	20,23,23	3.36	11 (55%)	26,34,34	3.39	10 (38%)
2	PO4	C	602	-	4,4,4	1.20	0	6,6,6	0.27	0
2	PO4	C	605	-	4,4,4	1.22	0	6,6,6	0.27	0
3	NMN	D	504	-	20,23,23	3.26	12 (60%)	26,34,34	3.43	10 (38%)
2	PO4	D	607	-	4,4,4	1.22	0	6,6,6	0.27	0
2	PO4	D	610	-	4,4,4	1.23	0	6,6,6	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NMN	E	505	-	20,23,23	3.32	10 (50%)	26,34,34	3.46	10 (38%)
2	PO4	E	608	-	4,4,4	1.04	0	6,6,6	0.27	0
2	PO4	E	609	-	4,4,4	1.14	0	6,6,6	0.27	0
3	NMN	F	506	-	20,23,23	3.34	10 (50%)	26,34,34	3.57	10 (38%)
2	PO4	F	611	-	4,4,4	1.03	0	6,6,6	0.27	0
2	PO4	F	612	-	4,4,4	1.19	0	6,6,6	0.27	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	NMN	A	501	-	-	0/10/30/30	0/2/2/2
2	PO4	A	601	-	-	0/0/0/0	0/0/0/0
2	PO4	A	606	-	-	0/0/0/0	0/0/0/0
3	NMN	B	502	-	-	0/10/30/30	0/2/2/2
2	PO4	B	603	-	-	0/0/0/0	0/0/0/0
2	PO4	B	604	-	-	0/0/0/0	0/0/0/0
3	NMN	C	503	-	-	0/10/30/30	0/2/2/2
2	PO4	C	602	-	-	0/0/0/0	0/0/0/0
2	PO4	C	605	-	-	0/0/0/0	0/0/0/0
3	NMN	D	504	-	-	0/10/30/30	0/2/2/2
2	PO4	D	607	-	-	0/0/0/0	0/0/0/0
2	PO4	D	610	-	-	0/0/0/0	0/0/0/0
3	NMN	E	505	-	-	0/10/30/30	0/2/2/2
2	PO4	E	608	-	-	0/0/0/0	0/0/0/0
2	PO4	E	609	-	-	0/0/0/0	0/0/0/0
3	NMN	F	506	-	-	0/10/30/30	0/2/2/2
2	PO4	F	611	-	-	0/0/0/0	0/0/0/0
2	PO4	F	612	-	-	0/0/0/0	0/0/0/0

All (63) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	NMN	C7-N7	-5.69	1.21	1.33
3	E	505	NMN	C7-N7	-5.68	1.21	1.33
3	C	503	NMN	C7-N7	-5.64	1.21	1.33
3	F	506	NMN	C7-N7	-5.53	1.21	1.33
3	D	504	NMN	C7-N7	-5.42	1.22	1.33
3	A	501	NMN	C7-N7	-5.35	1.22	1.33
3	F	506	NMN	C5R-C4R	-2.25	1.44	1.51

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	505	NMN	C5R-C4R	-2.16	1.44	1.51
3	B	502	NMN	C5R-C4R	-2.13	1.44	1.51
3	C	503	NMN	C5R-C4R	-2.08	1.44	1.51
3	D	504	NMN	C5R-C4R	-2.05	1.45	1.51
3	A	501	NMN	C5R-C4R	-2.04	1.45	1.51
3	D	504	NMN	P-O2P	-2.00	1.47	1.54
3	D	504	NMN	C2-C3	2.04	1.42	1.39
3	C	503	NMN	C2-C3	2.04	1.42	1.39
3	D	504	NMN	C6-C5	2.77	1.44	1.38
3	C	503	NMN	C6-C5	2.77	1.44	1.38
3	A	501	NMN	C6-C5	3.03	1.45	1.38
3	F	506	NMN	C6-C5	3.03	1.45	1.38
3	B	502	NMN	C6-C5	3.05	1.45	1.38
3	E	505	NMN	C6-C5	3.06	1.45	1.38
3	D	504	NMN	P-O3P	3.17	1.61	1.51
3	E	505	NMN	P-O3P	3.19	1.61	1.51
3	C	503	NMN	P-O3P	3.19	1.61	1.51
3	F	506	NMN	P-O3P	3.28	1.62	1.51
3	A	501	NMN	P-O3P	3.40	1.62	1.51
3	B	502	NMN	P-O3P	3.44	1.62	1.51
3	D	504	NMN	C5-C4	3.45	1.46	1.38
3	C	503	NMN	C5-C4	3.61	1.46	1.38
3	A	501	NMN	C5-C4	3.68	1.46	1.38
3	B	502	NMN	C5-C4	3.71	1.46	1.38
3	E	505	NMN	C5-C4	3.76	1.46	1.38
3	F	506	NMN	C5-C4	3.89	1.46	1.38
3	B	502	NMN	O4R-C4R	4.01	1.54	1.45
3	E	505	NMN	O4R-C4R	4.18	1.54	1.45
3	D	504	NMN	O4R-C4R	4.22	1.54	1.45
3	B	502	NMN	C3-C7	4.26	1.57	1.50
3	C	503	NMN	O4R-C4R	4.39	1.55	1.45
3	A	501	NMN	C3-C7	4.39	1.57	1.50
3	D	504	NMN	C3-C7	4.41	1.57	1.50
3	E	505	NMN	C3-C7	4.44	1.57	1.50
3	F	506	NMN	C3-C7	4.46	1.57	1.50
3	F	506	NMN	O4R-C4R	4.50	1.55	1.45
3	F	506	NMN	C6-N1	4.59	1.47	1.35
3	A	501	NMN	O4R-C4R	4.62	1.55	1.45
3	C	503	NMN	C3-C7	4.68	1.57	1.50
3	D	504	NMN	C6-N1	4.70	1.48	1.35
3	E	505	NMN	C6-N1	4.81	1.48	1.35
3	F	506	NMN	O7-C7	4.89	1.34	1.24

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	503	NMN	C6-N1	4.94	1.48	1.35
3	E	505	NMN	O7-C7	4.94	1.34	1.24
3	A	501	NMN	O7-C7	4.99	1.34	1.24
3	D	504	NMN	O7-C7	5.04	1.34	1.24
3	A	501	NMN	C6-N1	5.06	1.49	1.35
3	B	502	NMN	C6-N1	5.08	1.49	1.35
3	B	502	NMN	O7-C7	5.12	1.35	1.24
3	C	503	NMN	O7-C7	5.20	1.35	1.24
3	D	504	NMN	C4-C3	7.15	1.51	1.39
3	F	506	NMN	C4-C3	7.24	1.51	1.39
3	E	505	NMN	C4-C3	7.34	1.51	1.39
3	C	503	NMN	C4-C3	7.35	1.51	1.39
3	B	502	NMN	C4-C3	7.69	1.52	1.39
3	A	501	NMN	C4-C3	7.70	1.52	1.39

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	506	NMN	O7-C7-C3	-10.06	108.61	119.59
3	A	501	NMN	O7-C7-C3	-9.92	108.75	119.59
3	B	502	NMN	O7-C7-C3	-9.68	109.02	119.59
3	E	505	NMN	O7-C7-C3	-9.53	109.19	119.59
3	D	504	NMN	O7-C7-C3	-9.36	109.36	119.59
3	C	503	NMN	O7-C7-C3	-9.28	109.46	119.59
3	F	506	NMN	C5-C4-C3	-4.79	114.31	120.33
3	A	501	NMN	C5-C4-C3	-4.68	114.44	120.33
3	D	504	NMN	C4-C3-C7	-4.68	108.71	121.09
3	E	505	NMN	C5-C4-C3	-4.67	114.47	120.33
3	D	504	NMN	C5-C4-C3	-4.62	114.53	120.33
3	C	503	NMN	C4-C3-C7	-4.58	108.98	121.09
3	B	502	NMN	C5-C4-C3	-4.54	114.63	120.33
3	A	501	NMN	C4-C3-C7	-4.48	109.26	121.09
3	E	505	NMN	C4-C3-C7	-4.38	109.50	121.09
3	C	503	NMN	C5-C4-C3	-4.37	114.83	120.33
3	B	502	NMN	C4-C3-C7	-4.32	109.68	121.09
3	F	506	NMN	C4-C3-C7	-4.26	109.83	121.09
3	E	505	NMN	C5-C6-N1	-3.62	114.20	120.47
3	B	502	NMN	C5-C6-N1	-3.56	114.32	120.47
3	F	506	NMN	C5-C6-N1	-3.53	114.37	120.47
3	C	503	NMN	C5-C6-N1	-3.44	114.52	120.47
3	A	501	NMN	C5-C6-N1	-3.31	114.74	120.47
3	D	504	NMN	C5-C6-N1	-3.25	114.85	120.47

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	506	NMN	C4R-O4R-C1R	-3.24	106.16	109.72
3	A	501	NMN	C4R-O4R-C1R	-3.22	106.18	109.72
3	D	504	NMN	C4R-O4R-C1R	-3.15	106.26	109.72
3	C	503	NMN	C4R-O4R-C1R	-2.84	106.60	109.72
3	B	502	NMN	C4R-O4R-C1R	-2.42	107.06	109.72
3	E	505	NMN	C4R-O4R-C1R	-2.37	107.11	109.72
3	A	501	NMN	O4R-C1R-N1	2.35	110.71	108.13
3	D	504	NMN	O4R-C1R-N1	2.37	110.74	108.13
3	C	503	NMN	O4R-C1R-N1	2.39	110.76	108.13
3	B	502	NMN	O4R-C1R-N1	2.44	110.81	108.13
3	E	505	NMN	O4R-C1R-N1	2.92	111.34	108.13
3	B	502	NMN	C2-C3-C7	2.97	127.94	119.31
3	F	506	NMN	C2-C3-C7	2.98	127.96	119.31
3	A	501	NMN	C2-C3-C7	3.07	128.23	119.31
3	E	505	NMN	C2-C3-C7	3.09	128.27	119.31
3	C	503	NMN	C2-C3-C7	3.28	128.84	119.31
3	D	504	NMN	C2-C3-C7	3.29	128.87	119.31
3	F	506	NMN	O4R-C1R-N1	3.45	111.92	108.13
3	C	503	NMN	C2-C3-C4	3.50	122.18	118.29
3	F	506	NMN	C2-C3-C4	3.53	122.22	118.29
3	E	505	NMN	C2-C3-C4	3.53	122.22	118.29
3	B	502	NMN	C2-C3-C4	3.67	122.37	118.29
3	D	504	NMN	C2-C3-C4	3.71	122.42	118.29
3	A	501	NMN	C2-C3-C4	3.79	122.51	118.29
3	B	502	NMN	C6-C5-C4	4.02	125.51	119.44
3	A	501	NMN	C6-C5-C4	4.07	125.60	119.44
3	C	503	NMN	C6-C5-C4	4.16	125.73	119.44
3	F	506	NMN	C6-C5-C4	4.16	125.73	119.44
3	E	505	NMN	C6-C5-C4	4.19	125.78	119.44
3	D	504	NMN	C6-C5-C4	4.20	125.79	119.44
3	D	504	NMN	C3-C7-N7	9.48	128.20	117.82
3	C	503	NMN	C3-C7-N7	9.54	128.26	117.82
3	B	502	NMN	C3-C7-N7	9.62	128.34	117.82
3	E	505	NMN	C3-C7-N7	9.79	128.53	117.82
3	A	501	NMN	C3-C7-N7	9.80	128.54	117.82
3	F	506	NMN	C3-C7-N7	10.17	128.95	117.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 7 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	PO4	1	0
3	C	503	NMN	1	0
2	C	605	PO4	1	0
3	D	504	NMN	1	0
2	D	607	PO4	1	0
3	E	505	NMN	1	0
3	F	506	NMN	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	461/491 (93%)	0.19	14 (3%) 54 53	22, 39, 53, 67	0
1	B	464/491 (94%)	0.01	10 (2%) 65 64	19, 36, 53, 66	0
1	C	461/491 (93%)	0.40	24 (5%) 31 30	21, 42, 67, 81	0
1	D	464/491 (94%)	-0.33	7 (1%) 76 75	14, 23, 38, 69	0
1	E	464/491 (94%)	-0.32	1 (0%) 95 95	14, 24, 38, 57	0
1	F	463/491 (94%)	-0.41	5 (1%) 82 82	11, 20, 36, 64	0
All	All	2777/2946 (94%)	-0.08	61 (2%) 65 64	11, 30, 55, 81	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	484	ILE	7.7
1	A	484	ILE	4.8
1	B	338	SER	4.7
1	A	29	ASN	4.5
1	A	409	LEU	4.2
1	C	285	ASN	4.0
1	D	452	GLU	3.9
1	C	409	LEU	3.9
1	D	453	TYR	3.8
1	C	182	ASN	3.7
1	C	93	ASP	3.6
1	C	351	ILE	3.5
1	C	436	PRO	3.3
1	A	404	VAL	3.3
1	C	92	GLN	3.2
1	A	252	ALA	3.2
1	B	29	ASN	3.2
1	C	450	LEU	3.1
1	F	453	TYR	3.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	F	484	ILE	2.9
1	C	177	LEU	2.8
1	C	455	GLN	2.7
1	A	26	TYR	2.7
1	C	439	ASN	2.7
1	B	409	LEU	2.7
1	F	452	GLU	2.7
1	B	411	ILE	2.6
1	A	94	ASP	2.5
1	C	371	LYS	2.5
1	B	337	ASN	2.5
1	D	450	LEU	2.5
1	C	88	LYS	2.5
1	D	451	GLU	2.4
1	C	483	ASN	2.4
1	D	454	GLY	2.4
1	B	95	VAL	2.4
1	B	271	SER	2.3
1	A	370	GLN	2.3
1	A	420	ASP	2.3
1	D	271	SER	2.3
1	A	450	LEU	2.3
1	D	305	GLN	2.2
1	B	94	ASP	2.2
1	C	94	ASP	2.2
1	C	271	SER	2.2
1	C	411	ILE	2.2
1	E	484	ILE	2.2
1	F	451	GLU	2.1
1	C	328	LEU	2.1
1	A	419	ALA	2.1
1	C	181	GLY	2.1
1	A	181	GLY	2.1
1	B	484	ILE	2.1
1	F	450	LEU	2.0
1	C	255	LYS	2.0
1	C	309	ILE	2.0
1	C	244	ALA	2.0
1	B	483	ASN	2.0
1	A	244	ALA	2.0
1	A	245	ALA	2.0
1	C	343	LEU	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(Å <sup>2</sup> )	Q<0.9
2	PO4	E	608	5/5	0.59	0.54	10.31	87,88,89,90	0
3	NMN	E	505	22/22	0.94	0.19	2.42	17,31,42,46	0
3	NMN	D	504	22/22	0.94	0.19	1.74	16,26,31,33	0
3	NMN	F	506	22/22	0.95	0.17	1.16	9,28,34,37	0
3	NMN	B	502	22/22	0.93	0.20	1.09	26,38,41,47	0
3	NMN	A	501	22/22	0.93	0.19	0.80	27,37,40,42	0
3	NMN	C	503	22/22	0.94	0.18	0.43	36,39,41,43	0
2	PO4	C	602	5/5	0.98	0.17	0.06	47,48,49,50	0
2	PO4	B	603	5/5	0.99	0.15	-0.26	27,28,30,30	0
2	PO4	A	606	5/5	0.96	0.14	-0.33	49,51,52,52	0
2	PO4	D	610	5/5	0.98	0.12	-0.48	33,34,36,37	0
2	PO4	D	607	5/5	0.99	0.11	-1.24	15,15,17,19	0
2	PO4	B	604	5/5	0.98	0.10	-1.30	36,36,37,37	0
2	PO4	C	605	5/5	0.99	0.13	-1.43	35,36,38,40	0
2	PO4	F	612	5/5	0.99	0.08	-1.48	27,27,29,30	0
2	PO4	F	611	5/5	1.00	0.08	-1.59	16,17,19,20	0
2	PO4	A	601	5/5	0.96	0.10	-1.61	31,34,34,35	0
2	PO4	E	609	5/5	0.99	0.10	-1.72	22,22,23,24	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.