Full wwPDB X-ray Structure Validation Report (i)

Jul 13, 2021 – 10:03 AM EDT

PDB ID	:	1HZZ
Title	:	THE ASYMMETRIC COMPLEX OF THE TWO NUCLEOTIDE-BINDING COMPONENTS (DI, DIII) OF PROTON-TRANSLOCATING TRANSHYDROGENASE
Authors	:	Cotton, N.P.J.; White, S.A.; Peake, S.J.; McSweeney, S.; Jackson, J.B.
Deposited on	:	2001-01-27
Resolution	:	2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

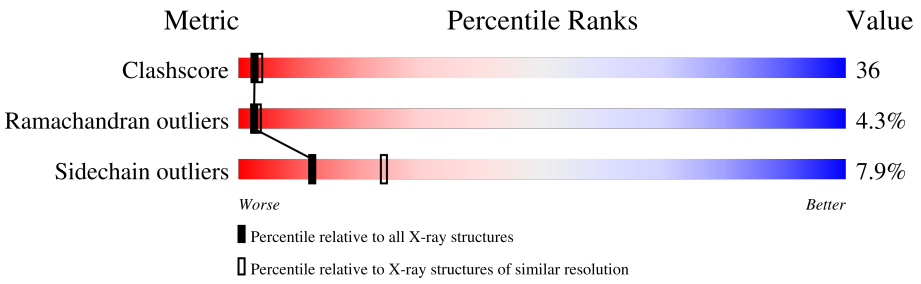
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.22

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	384	44% 42% 8% 6%
1	B	384	48% 38% 7% • 5%
2	C	203	34% 44% 7% 14%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6835 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.

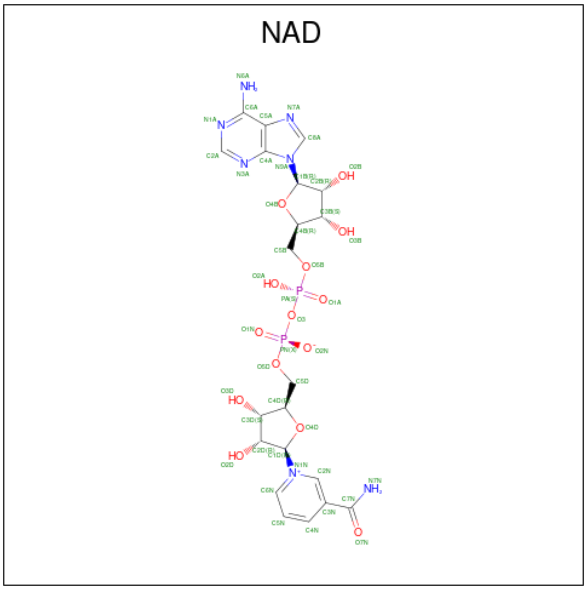
- Molecule 1 is a protein called PROTON-TRANSLOCATING NICOTINAMIDE NUCLEOTIDE TRANSHYDROGENASE SUBUNIT PNTAA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	362	2666	1684	460	505	17	0	0	0
1	B	366	2692	1701	465	509	17	0	0	0

- Molecule 2 is a protein called PROTON-TRANSLOCATING NICOTINAMIDE NUCLEOTIDE TRANSHYDROGENASE SUBUNIT PNTB.

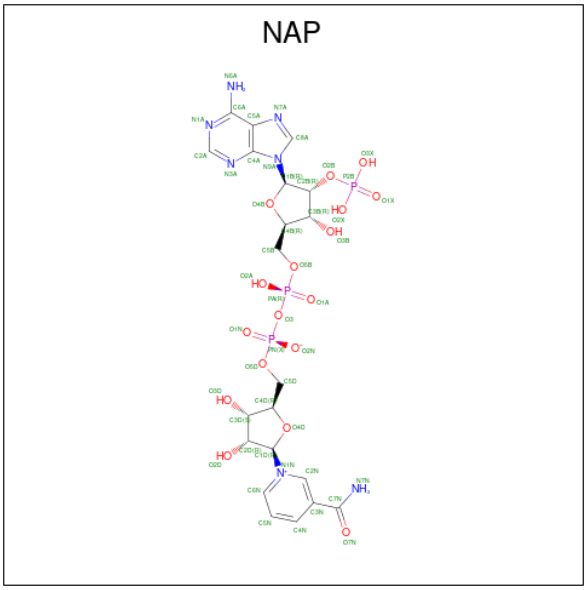
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	174	1311	830	217	253	11	0	0	0

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	C	1	48	21	7	17	3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	25	25	25	0	0
			Total	O		
5	B	40	40	40	0	0
			Total	O		
5	C	9	9	9	0	0

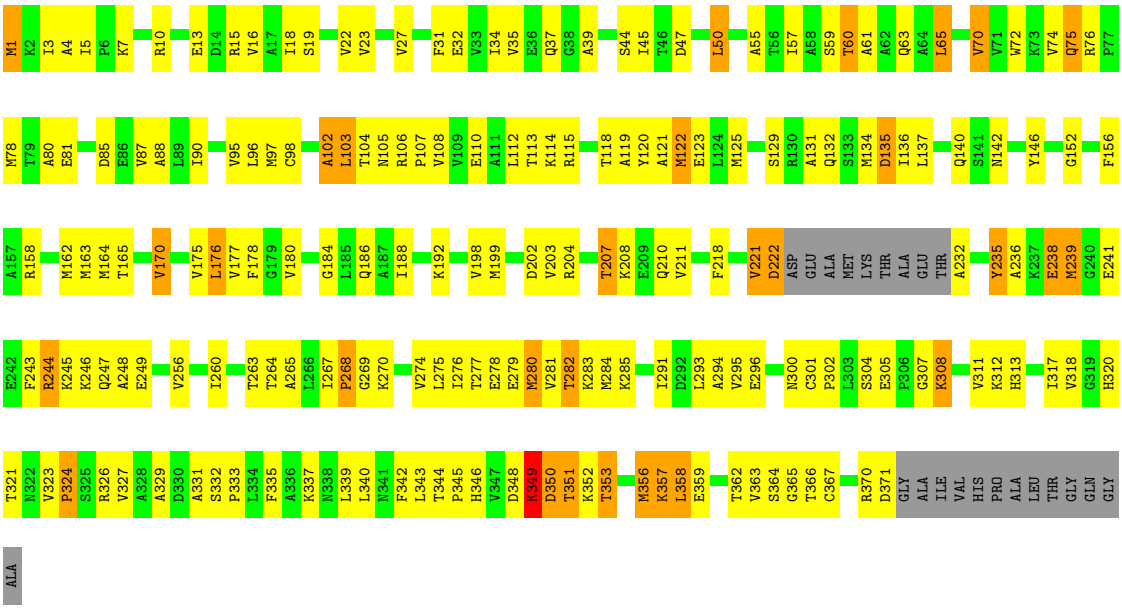
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

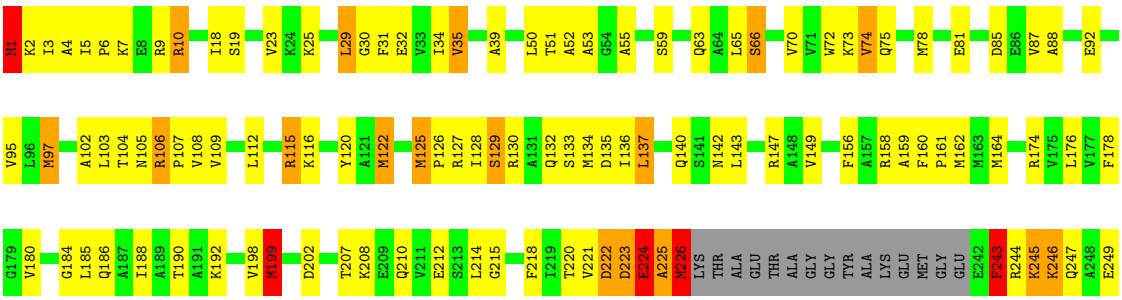
• Molecule 1: PROTON-TRANSLOCATING NICOTINAMIDE NUCLEOTIDE TRANSHYDROGENASE SUBUNIT PNTAA

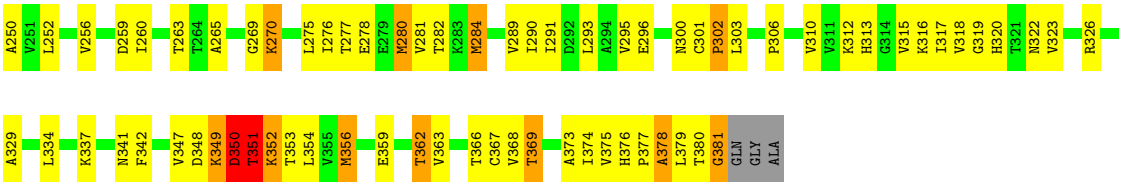
Chain A: 44% 42% 8% 6%



• Molecule 1: PROTON-TRANSLOCATING NICOTINAMIDE NUCLEOTIDE TRANSHYDROGENASE SUBUNIT PNTAA

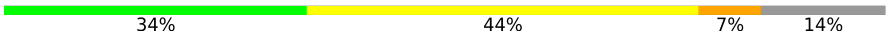
Chain B: 48% 38% 7% • 5%

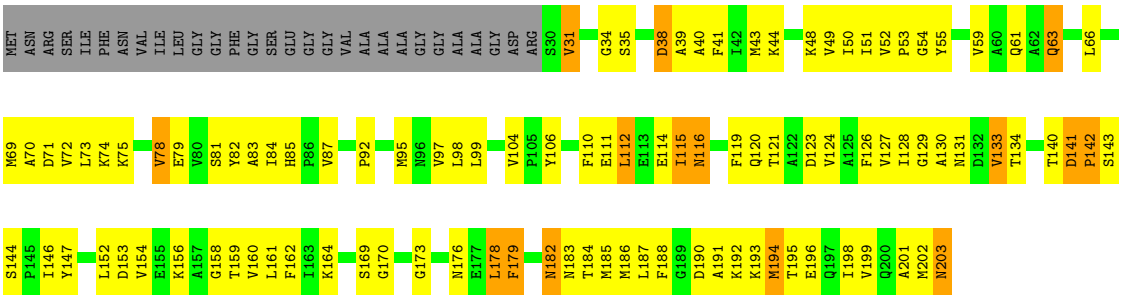




• Molecule 2: PROTON-TRANSLOCATING NICOTINAMIDE NUCLEOTIDE TRANSHYDROGENASE SUBUNIT PNTB

Chain C:





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	71.48Å 74.17Å 205.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.96 – 2.50	Depositor
% Data completeness (in resolution range)	95.0 (24.96-2.50)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.261 , 0.297	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6835	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

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: NAP, NAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	0.65	11/2700 (0.4%)	0.67	1/3656 (0.0%)
1	B	0.88	19/2727 (0.7%)	0.81	7/3697 (0.2%)
2	C	0.75	2/1334 (0.1%)	0.67	0/1803
All	All	0.77	32/6761 (0.5%)	0.73	8/9156 (0.1%)

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	381	GLY	C-O	25.74	1.64	1.23
2	C	203	ASN	C-OXT	21.29	1.63	1.23
1	B	226	MET	CG-SD	11.08	2.10	1.81
1	A	222	ASP	C-O	10.64	1.43	1.23
1	A	239	MET	CG-SD	8.28	2.02	1.81
1	B	348	ASP	C-N	7.13	1.50	1.34
2	C	203	ASN	C-O	6.74	1.36	1.23
1	B	224	GLU	CG-CD	6.61	1.61	1.51
1	B	1	MET	CG-SD	6.23	1.97	1.81
1	A	280	MET	CG-SD	6.04	1.96	1.81
1	A	284	MET	CG-SD	6.00	1.96	1.81
1	B	224	GLU	CD-OE2	5.97	1.32	1.25
1	A	134	MET	CG-SD	5.85	1.96	1.81
1	A	78	MET	CG-SD	5.78	1.96	1.81
1	B	223	ASP	C-O	5.76	1.34	1.23
1	A	125	MET	CG-SD	5.73	1.96	1.81
1	A	1	MET	CG-SD	5.72	1.96	1.81
1	B	226	MET	SD-CE	5.72	2.09	1.77
1	B	162	MET	CG-SD	5.65	1.95	1.81
1	B	356	MET	CG-SD	5.64	1.95	1.81
1	B	280	MET	CG-SD	5.61	1.95	1.81
1	A	356	MET	CG-SD	5.53	1.95	1.81
1	B	224	GLU	C-O	5.53	1.33	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	122	MET	CG-SD	5.42	1.95	1.81
1	B	284	MET	CG-SD	5.35	1.95	1.81
1	B	78	MET	CG-SD	5.33	1.95	1.81
1	A	162	MET	CG-SD	5.31	1.95	1.81
1	B	97	MET	CG-SD	5.26	1.94	1.81
1	B	164	MET	CG-SD	5.13	1.94	1.81
1	A	122	MET	CG-SD	5.04	1.94	1.81
1	B	199	MET	CG-SD	5.03	1.94	1.81
1	B	125	MET	CG-SD	5.02	1.94	1.81

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	350	ASP	O-C-N	12.18	142.18	122.70
1	B	350	ASP	CA-C-N	-8.97	97.46	117.20
1	B	226	MET	CB-CG-SD	-8.12	88.03	112.40
1	B	350	ASP	N-CA-C	-7.04	91.99	111.00
1	B	351	THR	O-C-N	6.45	133.02	122.70
1	B	351	THR	CA-C-N	-5.60	104.89	117.20
1	A	32	GLU	N-CA-C	-5.51	96.13	111.00
1	B	223	ASP	N-CA-C	-5.20	96.95	111.00

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	2666	0	2794	194	0
1	B	2692	0	2824	196	0
2	C	1311	0	1303	117	0
3	A	44	0	23	3	0
4	C	48	0	25	3	0
5	A	25	0	0	3	0
5	B	40	0	0	8	0
5	C	9	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6835	0	6969	501	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 36.

All (501) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:MET:SD	1:A:239:MET:CE	2.03	1.47
1:A:239:MET:SD	1:A:239:MET:CG	2.02	1.45
1:B:226:MET:CE	1:B:226:MET:SD	2.09	1.40
1:B:226:MET:SD	1:B:226:MET:CG	2.10	1.39
2:C:203:ASN:C	2:C:203:ASN:OXT	1.63	1.34
1:B:381:GLY:C	1:B:381:GLY:O	1.64	1.32
1:B:349:LYS:O	1:B:350:ASP:CG	1.86	1.14
1:B:351:THR:O	1:B:352:LYS:HG3	1.49	1.12
1:A:277:THR:H	1:A:280:MET:HE3	0.98	1.11
1:B:115:ARG:HH11	1:B:115:ARG:HB3	1.13	1.08
1:A:221:VAL:HG22	1:A:222:ASP:H	1.15	1.08
1:B:349:LYS:O	1:B:350:ASP:OD2	1.76	1.02
2:C:127:VAL:HG13	2:C:130:ALA:HB3	1.43	1.00
1:B:226:MET:SD	1:B:226:MET:CA	2.51	0.98
1:A:277:THR:N	1:A:280:MET:HE3	1.82	0.92
1:A:65:LEU:H	1:A:65:LEU:HD22	1.34	0.92
1:B:2:LYS:HE2	1:B:34:ILE:HD11	1.50	0.92
1:B:226:MET:HE2	5:B:421:HOH:O	1.68	0.92
2:C:53:PRO:HB2	2:C:95:MET:HE2	1.51	0.92
1:A:275:LEU:HD12	1:A:300:ASN:HB3	1.50	0.92
2:C:195:THR:O	2:C:199:VAL:HG23	1.72	0.90
2:C:111:GLU:HG3	2:C:114:GLU:HG2	1.54	0.89
1:B:226:MET:SD	1:B:226:MET:CB	2.60	0.89
1:B:10:ARG:HG3	1:B:10:ARG:HH11	1.40	0.87
1:A:277:THR:H	1:A:280:MET:CE	1.86	0.86
1:B:224:GLU:OE1	1:B:226:MET:N	2.09	0.85
1:A:72:TRP:HE1	1:A:97:MET:HE3	1.40	0.85
1:A:158:ARG:HG2	1:B:329:ALA:HB3	1.56	0.85
1:B:351:THR:O	1:B:352:LYS:CG	2.25	0.85
1:A:7:LYS:HG2	1:A:39:ALA:HA	1.58	0.83
1:B:351:THR:O	1:B:351:THR:HG23	1.78	0.83
2:C:99:LEU:HB3	2:C:104:VAL:HG11	1.61	0.82
2:C:99:LEU:HB3	2:C:104:VAL:CG1	2.09	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:VAL:HG22	1:A:222:ASP:N	1.96	0.81
1:B:368:VAL:HG12	1:B:369:THR:HG22	1.61	0.81
1:A:87:VAL:HG13	1:A:112:LEU:HD23	1.62	0.81
1:A:207:THR:HA	1:A:210:GLN:HE21	1.44	0.81
2:C:53:PRO:HB2	2:C:95:MET:CE	2.11	0.79
2:C:127:VAL:CG1	2:C:130:ALA:HB3	2.12	0.79
1:B:226:MET:SD	1:B:226:MET:HA	2.23	0.79
2:C:39:ALA:HB2	2:C:187:LEU:HD21	1.65	0.78
1:A:106:ARG:N	1:A:107:PRO:HD2	1.98	0.78
1:A:3:ILE:HG13	1:A:70:VAL:HG22	1.65	0.77
1:B:252:LEU:O	1:B:256:VAL:HG23	1.85	0.77
1:B:115:ARG:HH11	1:B:115:ARG:CB	1.94	0.76
1:B:87:VAL:HG23	1:B:112:LEU:HD23	1.68	0.76
1:B:221:VAL:HG11	1:B:247:GLN:HA	1.69	0.75
1:B:276:ILE:HG12	5:B:402:HOH:O	1.86	0.75
1:A:70:VAL:HB	1:A:95:VAL:HG22	1.69	0.74
2:C:66:LEU:HD13	2:C:128:ILE:HD13	1.70	0.73
1:B:25:LYS:O	1:B:29:LEU:HD12	1.87	0.73
2:C:85:HIS:CE1	2:C:87:VAL:HG22	2.23	0.73
1:B:115:ARG:HB3	1:B:115:ARG:NH1	1.96	0.73
2:C:50:ILE:HG22	2:C:81:SER:HB2	1.70	0.73
1:A:329:ALA:HB3	1:B:158:ARG:HG3	1.71	0.72
1:B:88:ALA:HA	1:B:115:ARG:HE	1.54	0.72
1:A:221:VAL:CG2	1:A:222:ASP:H	1.99	0.71
2:C:154:VAL:HG23	2:C:178:LEU:HD11	1.71	0.71
2:C:112:LEU:C	2:C:112:LEU:HD23	2.11	0.71
1:A:222:ASP:HA	1:A:246:LYS:HD2	1.71	0.70
1:B:222:ASP:OD2	1:B:225:ALA:HA	1.91	0.70
1:B:226:MET:SD	1:B:226:MET:N	2.64	0.70
1:B:178:PHE:HB3	1:B:275:LEU:HD13	1.73	0.70
1:B:97:MET:HB3	1:B:122:MET:CE	2.22	0.70
2:C:116:ASN:OD1	2:C:152:LEU:HA	1.93	0.69
2:C:69:MET:O	2:C:73:LEU:HB2	1.93	0.69
1:A:113:THR:HG23	1:A:371:ASP:OD2	1.93	0.68
1:A:296:GLU:HB2	5:A:420:HOH:O	1.94	0.68
1:A:244:ARG:HH11	1:A:244:ARG:HG2	1.59	0.68
1:B:120:TYR:OH	1:B:356:MET:HG3	1.94	0.68
2:C:194:MET:O	2:C:198:ILE:HG13	1.93	0.68
1:A:326:ARG:HG2	1:A:326:ARG:HH11	1.59	0.67
2:C:51:ILE:HG21	2:C:66:LEU:HD21	1.77	0.67
1:B:2:LYS:CE	1:B:34:ILE:HD11	2.23	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:SER:HB2	1:A:333:PRO:HD3	1.77	0.67
1:B:269:GLY:C	1:B:270:LYS:HD2	2.15	0.67
2:C:111:GLU:HB2	5:C:305:HOH:O	1.94	0.67
1:A:135:ASP:OD2	3:A:400:NAD:H4N	1.95	0.66
1:B:192:LYS:HE2	1:B:215:GLY:HA3	1.77	0.66
2:C:133:VAL:HG13	5:C:308:HOH:O	1.95	0.66
2:C:185:MET:O	2:C:186:MET:HE2	1.96	0.66
1:B:225:ALA:O	1:B:226:MET:C	2.34	0.65
1:A:244:ARG:HG2	1:A:244:ARG:NH1	2.10	0.65
1:B:374:ILE:HD12	1:B:374:ILE:H	1.61	0.65
1:A:221:VAL:HG23	1:A:247:GLN:HA	1.78	0.65
1:A:278:GLU:HA	1:A:281:VAL:HG22	1.78	0.65
1:B:222:ASP:CG	1:B:225:ALA:HA	2.17	0.65
1:A:47:ASP:HB3	1:A:57:ILE:CD1	2.27	0.65
1:B:129:SER:HA	1:B:132:GLN:HE21	1.61	0.65
1:A:275:LEU:HD12	1:A:300:ASN:CB	2.23	0.65
1:B:3:ILE:HG13	1:B:70:VAL:HG12	1.79	0.65
1:A:176:LEU:HD23	1:A:199:MET:HB3	1.77	0.65
2:C:146:ILE:HG12	2:C:146:ILE:O	1.97	0.65
2:C:161:LEU:HD12	2:C:185:MET:HB2	1.78	0.65
1:A:80:ALA:HB2	1:A:85:ASP:HB2	1.78	0.64
2:C:129:GLY:O	4:C:300:NAP:H4B	1.97	0.64
1:B:224:GLU:O	1:B:225:ALA:HB2	1.96	0.64
1:A:307:GLY:H	1:A:321:THR:HG23	1.61	0.64
1:B:226:MET:HE1	1:B:244:ARG:NH2	2.13	0.64
1:B:226:MET:CE	1:B:244:ARG:NH2	2.61	0.64
1:B:226:MET:CE	1:B:244:ARG:HH22	2.11	0.64
1:B:226:MET:HE2	1:B:244:ARG:HH22	1.61	0.64
2:C:85:HIS:HE1	2:C:87:VAL:HG22	1.63	0.64
1:A:158:ARG:HG2	1:B:329:ALA:CB	2.26	0.64
1:A:342:PHE:CE1	1:A:362:THR:HG22	2.32	0.63
1:A:178:PHE:CE2	1:A:276:ILE:HD11	2.32	0.63
2:C:159:THR:HA	2:C:183:ASN:HB3	1.81	0.63
2:C:140:THR:HG22	2:C:141:ASP:N	2.13	0.63
2:C:40:ALA:O	2:C:44:LYS:HB2	1.97	0.63
1:B:97:MET:HB3	1:B:122:MET:HE3	1.80	0.63
1:B:312:LYS:HG2	1:B:313:HIS:CD2	2.34	0.62
1:A:180:VAL:HG11	1:A:211:VAL:HG23	1.80	0.62
1:A:87:VAL:HG12	1:A:115:ARG:HG3	1.82	0.62
1:B:10:ARG:HH11	1:B:10:ARG:CG	2.09	0.62
1:B:224:GLU:CG	1:B:225:ALA:N	2.60	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:THR:O	1:B:351:THR:CG2	2.46	0.62
2:C:190:ASP:HB3	2:C:193:LYS:HD2	1.80	0.62
1:A:120:TYR:HB3	1:A:366:THR:HG22	1.82	0.62
1:B:174:ARG:HH21	1:B:199:MET:HE3	1.64	0.62
1:A:269:GLY:O	1:A:270:LYS:HG2	1.99	0.61
1:B:316:LYS:HG3	5:B:417:HOH:O	1.99	0.61
1:A:156:PHE:CZ	1:A:158:ARG:HB2	2.35	0.61
1:A:13:GLU:OE2	1:A:15:ARG:HB2	2.00	0.61
1:A:177:VAL:HG11	1:A:188:ILE:CG1	2.31	0.61
1:B:224:GLU:OE1	1:B:226:MET:SD	2.58	0.61
2:C:190:ASP:O	2:C:194:MET:HB2	2.01	0.60
1:A:307:GLY:N	1:A:321:THR:HG23	2.15	0.60
1:B:226:MET:HE1	1:B:244:ARG:CZ	2.31	0.60
1:B:85:ASP:OD2	1:B:88:ALA:HB2	2.02	0.60
2:C:63:GLN:HG3	2:C:98:LEU:HB3	1.84	0.60
2:C:124:VAL:HG22	2:C:159:THR:HG23	1.84	0.60
1:A:184:GLY:O	1:A:188:ILE:HG13	2.01	0.60
1:B:222:ASP:HA	1:B:224:GLU:HG3	1.82	0.60
1:A:105:ASN:O	1:A:108:VAL:HG12	2.02	0.60
1:B:6:PRO:HB3	1:B:65:LEU:HD11	1.83	0.60
1:B:18:ILE:HG13	1:B:19:SER:H	1.67	0.60
1:A:132:GLN:HG2	3:A:400:NAD:O7N	2.02	0.59
1:B:156:PHE:CE2	1:B:158:ARG:HB2	2.37	0.59
1:A:188:ILE:HG23	1:A:198:VAL:HG11	1.83	0.59
1:B:374:ILE:HD12	1:B:374:ILE:N	2.17	0.59
2:C:182:ASN:CG	2:C:183:ASN:N	2.55	0.59
1:A:178:PHE:HE2	1:A:276:ILE:HD11	1.68	0.59
1:A:106:ARG:N	1:A:107:PRO:CD	2.66	0.59
1:A:177:VAL:HG11	1:A:188:ILE:HG12	1.83	0.59
1:A:244:ARG:HH11	1:A:244:ARG:CG	2.15	0.59
1:A:264:THR:HG22	1:A:293:LEU:HD12	1.85	0.59
1:A:1:MET:N	1:A:352:LYS:NZ	2.51	0.58
2:C:55:TYR:O	2:C:59:VAL:HG23	2.03	0.58
2:C:141:ASP:C	2:C:143:SER:H	2.06	0.58
1:A:5:ILE:HD12	1:A:5:ILE:N	2.17	0.58
1:A:60:THR:HG23	1:A:63:GLN:H	1.68	0.58
1:A:110:GLU:O	1:A:114:LYS:HD3	2.03	0.58
2:C:169:SER:OG	2:C:173:GLY:HA2	2.04	0.58
1:A:343:LEU:C	1:A:343:LEU:HD23	2.23	0.58
1:B:221:VAL:CG1	1:B:247:GLN:HA	2.34	0.58
1:B:351:THR:O	1:B:352:LYS:CB	2.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:85:HIS:CG	2:C:133:VAL:HG22	2.38	0.58
1:B:224:GLU:OE1	1:B:225:ALA:N	2.37	0.57
2:C:170:GLY:HA3	4:C:300:NAP:O1A	2.04	0.57
1:A:23:VAL:O	1:A:27:VAL:HG23	2.04	0.57
1:A:22:VAL:HG22	1:A:340:LEU:HD22	1.87	0.57
1:B:142:ASN:HD21	1:B:186:GLN:HE21	1.50	0.57
1:B:376:HIS:CE1	1:B:378:ALA:HB3	2.38	0.57
1:B:224:GLU:OE2	1:B:226:MET:SD	2.62	0.57
1:A:87:VAL:CG1	1:A:112:LEU:HD23	2.33	0.57
2:C:192:LYS:O	2:C:196:GLU:HB2	2.05	0.57
1:B:379:LEU:O	1:B:381:GLY:N	2.38	0.57
2:C:99:LEU:O	2:C:104:VAL:HG12	2.05	0.57
2:C:124:VAL:HG22	2:C:159:THR:CG2	2.35	0.56
2:C:144:SER:OG	2:C:146:ILE:HG22	2.05	0.56
1:A:239:MET:SD	1:A:239:MET:CB	2.90	0.56
1:B:244:ARG:O	1:B:244:ARG:HG2	2.06	0.56
2:C:152:LEU:HB2	5:C:308:HOH:O	2.05	0.56
1:A:238:GLU:OE1	1:A:238:GLU:N	2.36	0.56
1:A:207:THR:HG22	1:A:218:PHE:CE1	2.40	0.56
1:A:70:VAL:HB	1:A:95:VAL:CG2	2.34	0.56
1:B:5:ILE:N	1:B:5:ILE:HD12	2.21	0.56
1:A:265:ALA:HB3	1:A:275:LEU:HD11	1.88	0.56
1:B:180:VAL:HG23	1:B:202:ASP:HB2	1.88	0.56
1:A:121:ALA:HB1	1:A:123:GLU:OE2	2.06	0.55
1:A:308:LYS:NZ	1:A:308:LYS:HB3	2.21	0.55
1:B:349:LYS:O	1:B:350:ASP:CB	2.52	0.55
1:A:1:MET:N	1:A:352:LYS:HZ3	2.05	0.55
1:B:160:PHE:N	1:B:161:PRO:HD2	2.21	0.55
2:C:182:ASN:CG	2:C:183:ASN:H	2.09	0.55
1:A:44:SER:HA	5:A:407:HOH:O	2.07	0.55
2:C:203:ASN:OXT	2:C:203:ASN:O	2.22	0.55
1:A:85:ASP:OD2	1:A:88:ALA:HB2	2.06	0.55
1:A:65:LEU:HD22	1:A:65:LEU:N	2.14	0.55
1:A:146:TYR:CE2	1:B:159:ALA:HA	2.41	0.55
1:A:279:GLU:O	1:A:282:THR:HG22	2.07	0.55
1:A:61:ALA:O	1:A:65:LEU:HD22	2.07	0.54
1:A:105:ASN:C	1:A:107:PRO:HD2	2.28	0.54
2:C:48:LYS:HD2	2:C:121:THR:O	2.05	0.54
1:B:51:THR:HA	1:B:55:ALA:O	2.06	0.54
1:B:322:ASN:HB3	5:B:389:HOH:O	2.06	0.54
1:B:282:THR:HG22	1:B:313:HIS:CE1	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:ALA:C	1:A:5:ILE:HD12	2.28	0.54
1:A:7:LYS:CG	1:A:39:ALA:HA	2.32	0.54
1:A:120:TYR:HE2	1:A:363:VAL:HG13	1.73	0.54
1:A:311:VAL:O	1:A:312:LYS:HG2	2.07	0.54
2:C:110:PHE:HB2	2:C:115:ILE:HD12	1.90	0.54
1:B:174:ARG:HH21	1:B:199:MET:CE	2.21	0.53
2:C:203:ASN:OXT	2:C:203:ASN:CB	2.56	0.53
1:B:180:VAL:HG12	1:B:180:VAL:O	2.08	0.53
1:B:178:PHE:CE1	1:B:276:ILE:HD11	2.44	0.53
1:A:3:ILE:HG22	1:A:31:PHE:HB3	1.89	0.53
1:A:178:PHE:HB2	1:A:263:THR:HA	1.89	0.53
1:A:1:MET:H3	1:A:352:LYS:NZ	2.06	0.53
1:B:349:LYS:C	1:B:351:THR:N	2.28	0.53
1:B:222:ASP:OD1	1:B:226:MET:HG2	2.09	0.53
1:B:350:ASP:C	1:B:352:LYS:N	2.62	0.53
1:A:47:ASP:HB3	1:A:57:ILE:HD13	1.90	0.53
1:B:275:LEU:HB2	1:B:300:ASN:HB3	1.89	0.53
1:B:350:ASP:HB2	1:B:353:THR:OG1	2.09	0.53
2:C:187:LEU:O	2:C:194:MET:HE3	2.08	0.53
1:A:97:MET:O	1:A:98:CYS:HB3	2.09	0.53
1:A:202:ASP:HB3	1:A:207:THR:HG21	1.89	0.53
1:B:349:LYS:O	1:B:349:LYS:HD2	2.09	0.53
1:A:313:HIS:N	1:A:313:HIS:CD2	2.77	0.52
1:B:108:VAL:HG13	1:B:109:VAL:N	2.24	0.52
1:A:346:HIS:HD2	1:A:357:LYS:H	1.56	0.52
1:B:128:ILE:HG13	1:B:128:ILE:O	2.09	0.52
1:B:208:LYS:HG2	1:B:212:GLU:OE2	2.09	0.52
2:C:85:HIS:CD2	2:C:112:LEU:HD12	2.44	0.52
1:A:95:VAL:HA	1:A:118:THR:O	2.09	0.52
1:A:114:LYS:HD2	1:A:114:LYS:N	2.23	0.52
1:A:367:CYS:SG	1:A:370:ARG:HB3	2.50	0.52
1:B:368:VAL:HG13	1:B:379:LEU:HD12	1.90	0.52
2:C:38:ASP:O	2:C:41:PHE:N	2.42	0.52
2:C:83:ALA:HB1	2:C:152:LEU:HD22	1.91	0.52
1:A:177:VAL:O	1:A:177:VAL:HG13	2.10	0.52
1:A:348:ASP:OD2	1:A:353:THR:HG23	2.09	0.52
1:A:281:VAL:HG23	1:A:282:THR:N	2.25	0.52
1:B:88:ALA:HA	1:B:115:ARG:NE	2.24	0.52
1:B:291:ILE:HD13	1:B:318:VAL:HB	1.92	0.52
2:C:40:ALA:HB3	5:C:301:HOH:O	2.09	0.52
2:C:61:GLN:HG3	2:C:63:GLN:HE22	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:203:ASN:OXT	2:C:203:ASN:CA	2.53	0.52
2:C:160:VAL:HG11	2:C:162:PHE:HE1	1.75	0.51
2:C:176:ASN:O	2:C:179:PHE:HB2	2.10	0.51
1:A:326:ARG:HG2	1:A:326:ARG:NH1	2.24	0.51
1:B:19:SER:O	1:B:23:VAL:HG23	2.10	0.51
1:A:294:ALA:HB3	1:A:300:ASN:OD1	2.11	0.51
1:B:350:ASP:O	1:B:353:THR:HG23	2.11	0.51
2:C:44:LYS:HE2	2:C:202:MET:O	2.10	0.51
2:C:140:THR:O	2:C:142:PRO:HD3	2.11	0.51
2:C:142:PRO:HA	2:C:147:TYR:CD2	2.45	0.51
2:C:34:GLY:HA3	2:C:185:MET:CE	2.40	0.51
2:C:119:PHE:CD2	2:C:154:VAL:HA	2.45	0.51
1:A:10:ARG:HG3	5:A:417:HOH:O	2.10	0.51
2:C:82:TYR:CE1	2:C:104:VAL:HG21	2.46	0.51
2:C:84:ILE:HD11	2:C:99:LEU:HD11	1.92	0.51
1:A:65:LEU:H	1:A:65:LEU:CD2	2.14	0.51
1:B:97:MET:HB3	1:B:122:MET:HE1	1.90	0.51
1:B:374:ILE:H	1:B:374:ILE:CD1	2.21	0.51
1:A:163:MET:HB2	1:A:170:VAL:HG13	1.92	0.51
1:A:5:ILE:HG12	1:A:18:ILE:HB	1.93	0.51
1:A:274:VAL:HG13	1:A:274:VAL:O	2.10	0.50
2:C:78:VAL:HG12	2:C:79:GLU:N	2.25	0.50
1:B:277:THR:OG1	1:B:280:MET:HG3	2.11	0.50
1:A:97:MET:CE	1:A:343:LEU:HD12	2.42	0.50
1:A:3:ILE:CG1	1:A:70:VAL:HG22	2.37	0.50
1:A:106:ARG:O	1:A:106:ARG:HD3	2.12	0.50
1:A:129:SER:C	1:A:131:ALA:H	2.13	0.50
1:A:269:GLY:C	1:A:270:LYS:HG2	2.31	0.50
1:B:263:THR:HG21	1:B:276:ILE:HD13	1.94	0.50
1:B:120:TYR:HD1	1:B:366:THR:HG23	1.76	0.50
1:B:243:PHE:C	1:B:245:LYS:H	2.13	0.50
2:C:158:GLY:O	2:C:183:ASN:OD1	2.30	0.50
1:A:18:ILE:HG13	1:A:19:SER:N	2.27	0.50
1:A:16:VAL:HG22	1:A:18:ILE:HG22	1.94	0.50
1:B:31:PHE:HE1	1:B:347:VAL:HG21	1.77	0.50
1:A:165:THR:CG2	2:C:92:PRO:HB3	2.42	0.49
1:A:45:ILE:CG2	1:A:50:LEU:HD13	2.42	0.49
1:B:106:ARG:N	1:B:107:PRO:CD	2.75	0.49
1:A:323:VAL:HB	1:A:324:PRO:HD3	1.93	0.49
1:B:246:LYS:HE2	1:B:249:GLU:HG2	1.94	0.49
2:C:48:LYS:HA	2:C:79:GLU:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:MET:O	1:A:358:LEU:N	2.45	0.49
1:B:160:PHE:CZ	1:B:260:ILE:HD12	2.46	0.49
1:B:226:MET:HE2	1:B:244:ARG:NH2	2.28	0.49
1:B:320:HIS:HA	5:B:411:HOH:O	2.11	0.49
2:C:69:MET:HE1	2:C:126:PHE:CD2	2.48	0.49
2:C:191:ALA:HB2	4:C:300:NAP:H2A	1.94	0.49
1:A:291:ILE:HD13	1:A:318:VAL:HB	1.94	0.49
1:B:192:LYS:HG3	5:B:420:HOH:O	2.13	0.49
2:C:112:LEU:HD23	2:C:112:LEU:O	2.12	0.49
2:C:119:PHE:CE2	2:C:154:VAL:HG12	2.48	0.49
1:B:3:ILE:HA	1:B:70:VAL:O	2.13	0.49
2:C:66:LEU:HD13	2:C:128:ILE:HG21	1.94	0.49
2:C:185:MET:HB3	2:C:187:LEU:CD1	2.43	0.49
1:A:256:VAL:O	1:A:285:LYS:HG3	2.13	0.48
1:B:30:GLY:CA	1:B:352:LYS:HE2	2.43	0.48
1:B:224:GLU:CD	1:B:226:MET:SD	2.92	0.48
1:B:376:HIS:HE1	1:B:378:ALA:HB3	1.77	0.48
2:C:70:ALA:O	2:C:74:LYS:HG2	2.13	0.48
1:B:284:MET:CE	1:B:290:ILE:HD11	2.43	0.48
1:A:35:VAL:HG11	1:A:50:LEU:HD23	1.94	0.48
1:B:188:ILE:HG23	1:B:198:VAL:HG11	1.94	0.48
1:B:293:LEU:CD2	1:B:323:VAL:HG21	2.44	0.48
1:A:135:ASP:OD1	1:A:137:LEU:HB2	2.13	0.48
1:A:152:GLY:HA2	1:A:291:ILE:HD11	1.96	0.48
2:C:49:VAL:HG22	2:C:124:VAL:HB	1.96	0.48
1:A:23:VAL:HG11	1:A:55:ALA:HB2	1.95	0.48
1:A:3:ILE:HA	1:A:70:VAL:O	2.13	0.48
1:A:118:THR:HG23	1:A:370:ARG:HA	1.96	0.48
1:B:269:GLY:O	1:B:270:LYS:HD2	2.13	0.48
2:C:154:VAL:HG23	2:C:178:LEU:CD1	2.41	0.48
1:A:4:ALA:HA	1:A:34:ILE:O	2.14	0.48
1:A:357:LYS:O	1:A:359:GLU:N	2.44	0.47
1:A:120:TYR:OH	1:A:356:MET:HG3	2.14	0.47
2:C:160:VAL:HG11	2:C:162:PHE:CE1	2.48	0.47
1:B:1:MET:SD	1:B:354:LEU:HD22	2.54	0.47
1:A:207:THR:CG2	1:A:218:PHE:HE1	2.27	0.47
1:A:305:GLU:OE1	1:A:308:LYS:HD2	2.14	0.47
1:B:367:CYS:O	1:B:376:HIS:HB2	2.15	0.47
1:B:10:ARG:HG3	1:B:10:ARG:NH1	2.19	0.47
1:B:35:VAL:HG21	1:B:50:LEU:CD2	2.45	0.47
1:A:202:ASP:CB	1:A:207:THR:HG21	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:THR:HG22	1:B:363:VAL:HG23	1.97	0.47
1:A:75:GLN:HE21	1:A:75:GLN:HB2	1.51	0.47
1:B:342:PHE:CZ	1:B:366:THR:HG21	2.50	0.47
1:A:102:ALA:O	1:A:104:THR:N	2.48	0.47
1:B:221:VAL:HG22	1:B:247:GLN:HG2	1.97	0.47
2:C:134:THR:HG21	2:C:154:VAL:HG11	1.97	0.47
1:A:142:ASN:O	1:A:186:GLN:HG2	2.14	0.46
1:A:349:LYS:HE3	1:A:349:LYS:HA	1.97	0.46
1:B:35:VAL:HG21	1:B:50:LEU:HD23	1.97	0.46
2:C:35:SER:O	2:C:38:ASP:HB2	2.16	0.46
2:C:66:LEU:CD1	2:C:128:ILE:HG21	2.45	0.46
2:C:141:ASP:O	2:C:143:SER:N	2.47	0.46
1:A:265:ALA:HA	3:A:400:NAD:O4B	2.15	0.46
1:B:6:PRO:HD2	1:B:72:TRP:O	2.16	0.46
1:B:222:ASP:OD1	1:B:225:ALA:HA	2.15	0.46
1:B:301:CYS:HA	1:B:302:PRO:HD2	1.78	0.46
1:A:59:SER:HB2	1:A:63:GLN:NE2	2.30	0.46
1:A:244:ARG:HD2	1:A:274:VAL:CG1	2.46	0.46
1:B:221:VAL:CG2	1:B:244:ARG:HH21	2.28	0.46
1:B:373:ALA:O	1:B:375:VAL:HG23	2.16	0.46
1:A:72:TRP:NE1	1:A:97:MET:HE3	2.20	0.46
2:C:31:VAL:HG13	2:C:186:MET:HG3	1.96	0.46
1:A:327:VAL:O	1:A:327:VAL:HG12	2.16	0.46
2:C:52:VAL:HG12	2:C:52:VAL:O	2.15	0.46
1:B:178:PHE:CZ	1:B:276:ILE:HD11	2.51	0.46
2:C:85:HIS:CD2	2:C:133:VAL:HG22	2.50	0.46
2:C:169:SER:HB2	5:C:302:HOH:O	2.15	0.46
1:A:72:TRP:CD2	1:A:339:LEU:HD13	2.51	0.46
1:B:224:GLU:CG	1:B:225:ALA:H	2.28	0.46
1:B:315:VAL:O	1:B:317:ILE:HD12	2.15	0.46
2:C:142:PRO:HB3	2:C:147:TYR:CZ	2.51	0.46
1:A:256:VAL:HG12	1:A:283:LYS:O	2.15	0.45
2:C:162:PHE:CD2	2:C:179:PHE:HE1	2.34	0.45
1:A:60:THR:HG23	1:A:63:GLN:HB2	1.99	0.45
1:A:72:TRP:CG	1:A:339:LEU:HD13	2.52	0.45
1:B:102:ALA:C	1:B:104:THR:H	2.19	0.45
1:A:140:GLN:OE1	1:A:331:ALA:HB1	2.16	0.45
1:B:7:LYS:HG3	1:B:39:ALA:HA	1.98	0.45
1:B:128:ILE:O	1:B:130:ARG:N	2.50	0.45
1:B:149:VAL:HG21	1:B:190:THR:HB	1.98	0.45
1:B:51:THR:C	1:B:53:ALA:H	2.20	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ALA:HB3	1:B:300:ASN:ND2	2.32	0.45
1:B:320:HIS:HB3	5:B:410:HOH:O	2.17	0.45
2:C:63:GLN:CG	2:C:98:LEU:HB3	2.46	0.45
1:A:106:ARG:NH1	1:A:110:GLU:HB3	2.32	0.45
1:A:239:MET:HB3	1:A:243:PHE:CG	2.52	0.45
1:B:156:PHE:CZ	1:B:259:ASP:HB3	2.52	0.45
1:B:323:VAL:O	1:B:326:ARG:HB2	2.16	0.45
2:C:123:ASP:O	2:C:159:THR:HG22	2.17	0.45
2:C:187:LEU:C	2:C:194:MET:HE3	2.36	0.45
1:B:295:VAL:HG11	1:B:319:GLY:O	2.16	0.45
1:A:136:ILE:O	1:A:140:GLN:HG2	2.17	0.44
1:B:70:VAL:HG22	1:B:95:VAL:HB	1.99	0.44
2:C:164:LYS:O	2:C:188:PHE:HA	2.18	0.44
1:A:97:MET:HE3	1:A:343:LEU:HD12	1.98	0.44
1:A:344:THR:HB	1:A:345:PRO:HD3	2.00	0.44
1:B:143:LEU:HD11	1:B:334:LEU:HD12	1.98	0.44
1:A:308:LYS:HB3	1:A:308:LYS:HZ3	1.81	0.44
1:B:246:LYS:HD2	1:B:246:LYS:HA	1.79	0.44
2:C:140:THR:HG22	2:C:141:ASP:H	1.79	0.44
1:A:1:MET:H3	1:A:352:LYS:HZ1	1.64	0.44
2:C:43:MET:HE2	2:C:73:LEU:HD21	2.00	0.44
1:A:178:PHE:CD2	1:A:263:THR:HG22	2.52	0.44
1:A:337:LYS:HE3	1:A:337:LYS:HB2	1.83	0.44
1:B:10:ARG:CG	1:B:10:ARG:NH1	2.72	0.44
1:B:105:ASN:O	1:B:108:VAL:HG12	2.18	0.44
1:B:208:LYS:HG3	1:B:218:PHE:CB	2.47	0.44
1:B:221:VAL:HG12	1:B:250:ALA:HB2	2.00	0.44
1:B:224:GLU:HG3	1:B:225:ALA:N	2.33	0.44
1:B:282:THR:HG22	1:B:313:HIS:ND1	2.33	0.44
1:B:295:VAL:CG1	1:B:296:GLU:N	2.80	0.44
1:A:122:MET:HB2	1:A:136:ILE:HD13	1.99	0.44
1:A:156:PHE:CE2	1:A:158:ARG:HB2	2.53	0.44
1:B:246:LYS:HE2	1:B:249:GLU:CG	2.47	0.44
1:A:276:ILE:O	1:A:302:PRO:HD2	2.18	0.44
1:B:349:LYS:C	1:B:349:LYS:HD2	2.38	0.44
1:A:180:VAL:HG11	1:A:211:VAL:CG2	2.46	0.43
1:B:122:MET:HB3	1:B:136:ILE:HD13	1.99	0.43
2:C:34:GLY:HA3	2:C:185:MET:HE3	2.00	0.43
2:C:72:VAL:HA	2:C:75:LYS:HB2	1.99	0.43
1:A:203:VAL:HG11	1:A:239:MET:HG3	1.99	0.43
1:A:204:ARG:O	1:A:207:THR:HB	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:ALA:HA	1:B:34:ILE:O	2.18	0.43
1:A:207:THR:HG22	1:A:208:LYS:N	2.33	0.43
1:A:305:GLU:O	1:A:308:LYS:HG2	2.18	0.43
1:A:332:SER:O	1:A:333:PRO:C	2.56	0.43
1:B:317:ILE:HD12	1:B:317:ILE:N	2.34	0.43
1:A:61:ALA:O	1:A:65:LEU:CD2	2.66	0.43
1:B:256:VAL:HG12	1:B:256:VAL:O	2.18	0.43
2:C:187:LEU:HB3	2:C:194:MET:CE	2.49	0.43
2:C:203:ASN:OXT	2:C:203:ASN:HB2	2.19	0.43
1:A:364:SER:OG	1:A:365:GLY:N	2.51	0.43
1:B:65:LEU:O	1:B:66:SER:C	2.57	0.43
1:B:220:THR:HB	5:B:421:HOH:O	2.18	0.43
2:C:78:VAL:CG1	2:C:79:GLU:N	2.81	0.43
1:A:35:VAL:HG11	1:A:50:LEU:CD2	2.49	0.43
2:C:97:VAL:HG13	2:C:98:LEU:N	2.33	0.43
2:C:140:THR:CG2	2:C:141:ASP:H	2.32	0.43
1:B:50:LEU:HD12	1:B:50:LEU:HA	1.83	0.43
1:B:51:THR:C	1:B:53:ALA:N	2.72	0.43
1:B:184:GLY:O	1:B:188:ILE:HG13	2.18	0.43
2:C:54:GLY:N	2:C:95:MET:HE1	2.34	0.43
2:C:119:PHE:CG	2:C:153:ASP:O	2.71	0.43
1:A:221:VAL:O	1:A:222:ASP:HB2	2.18	0.43
1:B:30:GLY:O	1:B:352:LYS:HD3	2.19	0.43
1:B:243:PHE:CE1	1:B:245:LYS:HE3	2.53	0.43
1:A:18:ILE:HG13	1:A:19:SER:H	1.84	0.43
1:A:74:VAL:HA	1:A:335:PHE:CE2	2.54	0.43
1:A:96:LEU:O	1:A:119:ALA:HA	2.19	0.43
1:A:312:LYS:HB3	1:A:313:HIS:CD2	2.54	0.43
1:A:348:ASP:O	1:A:350:ASP:N	2.52	0.42
1:B:7:LYS:HD3	1:B:9:ARG:NH1	2.34	0.42
1:B:136:ILE:O	1:B:140:GLN:HG2	2.19	0.42
2:C:112:LEU:C	2:C:112:LEU:CD2	2.82	0.42
2:C:161:LEU:HD12	2:C:161:LEU:HA	1.78	0.42
1:A:10:ARG:NH2	1:A:76:ARG:HH22	2.16	0.42
1:A:221:VAL:CG2	1:A:222:ASP:N	2.67	0.42
1:A:235:TYR:HB2	1:A:236:ALA:H	1.70	0.42
1:B:295:VAL:HG13	1:B:296:GLU:N	2.33	0.42
1:B:92:GLU:OE2	1:B:116:LYS:HD2	2.20	0.42
1:A:351:THR:O	1:A:353:THR:HG22	2.19	0.42
1:B:18:ILE:HG13	1:B:19:SER:N	2.32	0.42
1:A:329:ALA:HB3	1:B:158:ARG:CG	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:VAL:CG1	1:B:109:VAL:N	2.82	0.42
1:B:133:SER:HB2	1:B:341:ASN:ND2	2.34	0.42
1:A:10:ARG:NH2	1:A:76:ARG:NH2	2.67	0.42
1:B:135:ASP:OD1	1:B:137:LEU:HB2	2.19	0.42
1:B:125:MET:HA	1:B:126:PRO:HD3	1.90	0.42
1:B:185:LEU:HD23	1:B:185:LEU:HA	1.93	0.42
2:C:178:LEU:O	2:C:184:THR:HG21	2.19	0.42
2:C:187:LEU:CD1	2:C:187:LEU:N	2.83	0.42
2:C:40:ALA:HB1	2:C:201:ALA:O	2.20	0.42
2:C:119:PHE:CB	2:C:153:ASP:O	2.68	0.42
1:A:279:GLU:OE2	1:A:283:LYS:HE3	2.20	0.41
1:A:295:VAL:HG23	1:A:304:SER:OG	2.19	0.41
1:A:317:ILE:N	1:A:317:ILE:HD12	2.35	0.41
1:B:59:SER:HB2	1:B:63:GLN:OE1	2.19	0.41
1:B:281:VAL:HA	1:B:284:MET:CE	2.50	0.41
1:B:359:GLU:OE1	1:B:359:GLU:N	2.47	0.41
2:C:187:LEU:N	2:C:187:LEU:HD12	2.35	0.41
1:A:103:LEU:CD1	1:A:103:LEU:H	2.33	0.41
1:B:275:LEU:HD23	1:B:275:LEU:HA	1.87	0.41
1:A:192:LYS:HG2	1:A:198:VAL:CG2	2.51	0.41
1:B:160:PHE:N	1:B:161:PRO:CD	2.83	0.41
1:B:202:ASP:HB3	1:B:207:THR:HG21	2.02	0.41
2:C:71:ASP:HA	2:C:74:LYS:HG2	2.03	0.41
1:A:247:GLN:C	1:A:249:GLU:N	2.73	0.41
1:A:265:ALA:CB	1:A:275:LEU:HD11	2.50	0.41
1:B:81:GLU:O	1:B:81:GLU:HG3	2.20	0.41
1:B:214:LEU:HD23	1:B:214:LEU:HA	1.91	0.41
1:B:278:GLU:HG3	1:B:303:LEU:HD11	2.01	0.41
1:A:267:ILE:O	1:A:268:PRO:C	2.58	0.41
1:B:74:VAL:CG2	1:B:75:GLN:N	2.83	0.41
1:B:192:LYS:CE	1:B:215:GLY:HA3	2.47	0.41
1:B:303:LEU:O	1:B:310:VAL:HG11	2.21	0.41
1:B:350:ASP:C	1:B:352:LYS:H	2.23	0.41
2:C:53:PRO:HB2	2:C:95:MET:HE1	1.99	0.41
1:A:16:VAL:O	1:A:16:VAL:HG13	2.20	0.41
1:A:97:MET:HB3	1:A:122:MET:HE2	2.01	0.41
1:A:342:PHE:CE2	1:A:356:MET:HE1	2.54	0.41
1:B:109:VAL:HG13	1:B:368:VAL:CG1	2.51	0.41
1:A:103:LEU:H	1:A:103:LEU:HD13	1.86	0.41
1:A:176:LEU:HD13	1:A:178:PHE:CZ	2.55	0.41
1:A:177:VAL:HG11	1:A:188:ILE:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ALA:HB3	1:A:300:ASN:ND2	2.35	0.41
1:B:51:THR:O	1:B:53:ALA:N	2.54	0.41
2:C:61:GLN:HA	2:C:63:GLN:OE1	2.20	0.41
2:C:120:GLN:HG2	2:C:156:LYS:HD3	2.02	0.41
2:C:160:VAL:HB	2:C:184:THR:HA	2.03	0.41
1:B:18:ILE:HD12	1:B:18:ILE:HA	1.86	0.41
1:B:125:MET:CE	1:B:134:MET:HB3	2.51	0.41
1:B:342:PHE:CE1	1:B:362:THR:HG23	2.56	0.41
1:A:175:VAL:HG22	1:A:260:ILE:HB	2.03	0.40
1:A:232:ALA:HA	1:A:235:TYR:O	2.21	0.40
1:B:9:ARG:HD3	1:B:9:ARG:HA	1.83	0.40
1:B:104:THR:HG23	1:B:105:ASN:ND2	2.35	0.40
2:C:152:LEU:CB	5:C:308:HOH:O	2.68	0.40
1:A:203:VAL:HG12	1:A:243:PHE:CZ	2.56	0.40
1:B:275:LEU:C	1:B:276:ILE:HD12	2.41	0.40
1:A:241:GLU:OE2	1:A:245:LYS:NZ	2.44	0.40
1:A:352:LYS:HE2	1:A:352:LYS:HB3	1.90	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	358/384 (93%)	295 (82%)	50 (14%)	13 (4%)	3 4
1	B	362/384 (94%)	317 (88%)	29 (8%)	16 (4%)	2 3
2	C	172/203 (85%)	138 (80%)	25 (14%)	9 (5%)	2 2
All	All	892/971 (92%)	750 (84%)	104 (12%)	38 (4%)	2 3

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	102	ALA
1	A	103	LEU
1	A	349	LYS
1	A	357	LYS
1	B	66	SER
1	B	129	SER
1	B	224	GLU
1	B	225	ALA
1	B	351	THR
1	B	352	LYS
1	B	380	THR
2	C	182	ASN
1	A	221	VAL
1	B	243	PHE
1	B	378	ALA
2	C	78	VAL
2	C	131	ASN
1	A	238	GLU
1	A	350	ASP
1	B	245	LYS
2	C	31	VAL
2	C	38	ASP
2	C	141	ASP
1	A	268	PRO
1	A	320	HIS
1	A	351	THR
1	A	358	LEU
1	B	103	LEU
1	B	306	PRO
2	C	112	LEU
2	C	115	ILE
1	B	350	ASP
1	B	377	PRO
2	C	142	PRO
1	A	90	ILE
1	A	248	ALA
1	B	52	ALA
1	B	302	PRO

5.3.2 Protein sidechains (i)

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	282/296 (95%)	262 (93%)	20 (7%)	14	28
1	B	286/296 (97%)	257 (90%)	29 (10%)	7	14
2	C	138/154 (90%)	131 (95%)	7 (5%)	24	45
All	All	706/746 (95%)	650 (92%)	56 (8%)	12	24

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	50	LEU
1	A	60	THR
1	A	65	LEU
1	A	70	VAL
1	A	75	GLN
1	A	81	GLU
1	A	135	ASP
1	A	164	MET
1	A	170	VAL
1	A	176	LEU
1	A	207	THR
1	A	235	TYR
1	A	244	ARG
1	A	282	THR
1	A	301	CYS
1	A	308	LYS
1	A	324	PRO
1	A	349	LYS
1	A	353	THR
1	B	1	MET
1	B	10	ARG
1	B	29	LEU
1	B	32	GLU
1	B	35	VAL
1	B	73	LYS
1	B	74	VAL
1	B	106	ARG
1	B	115	ARG
1	B	127	ARG

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Mol	Chain	Res	Type
1	B	137	LEU
1	B	147	ARG
1	B	176	LEU
1	B	199	MET
1	B	210	GLN
1	B	222	ASP
1	B	223	ASP
1	B	224	GLU
1	B	226	MET
1	B	243	PHE
1	B	246	LYS
1	B	270	LYS
1	B	289	VAL
1	B	337	LYS
1	B	349	LYS
1	B	350	ASP
1	B	351	THR
1	B	362	THR
1	B	369	THR
2	C	63	GLN
2	C	106	TYR
2	C	116	ASN
2	C	133	VAL
2	C	178	LEU
2	C	179	PHE
2	C	194	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	63	GLN
1	A	75	GLN
1	A	210	GLN
1	A	247	GLN
1	A	313	HIS
1	A	338	ASN
1	A	346	HIS
1	B	99	HIS
1	B	105	ASN
1	B	132	GLN
1	B	142	ASN

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Mol	Chain	Res	Type
1	B	186	GLN
1	B	376	HIS
2	C	120	GLN
2	C	183	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	NAD	A	400	-	42,48,48	3.11	10 (23%)	50,73,73	2.02	12 (24%)
4	NAP	C	300	-	45,52,52	1.79	9 (20%)	56,80,80	1.86	14 (25%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAD	A	400	-	-	12/26/62/62	0/5/5/5
4	NAP	C	300	-	-	11/31/67/67	0/5/5/5

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	NAD	O2D-C2D	-12.61	1.13	1.43
3	A	400	NAD	C5D-C4D	-7.58	1.28	1.51
4	C	300	NAP	C2N-N1N	6.90	1.43	1.35
3	A	400	NAD	C3N-C7N	5.97	1.59	1.50
3	A	400	NAD	C6N-N1N	5.93	1.49	1.35
3	A	400	NAD	C2N-N1N	5.22	1.41	1.35
3	A	400	NAD	C4N-C3N	4.92	1.47	1.39
4	C	300	NAP	P2B-O2B	3.71	1.66	1.59
3	A	400	NAD	C4A-N3A	3.62	1.40	1.35
3	A	400	NAD	C2A-N1A	3.58	1.40	1.33
4	C	300	NAP	O4D-C1D	3.36	1.45	1.41
4	C	300	NAP	C6N-N1N	3.16	1.43	1.35
3	A	400	NAD	O4D-C1D	3.14	1.45	1.41
3	A	400	NAD	PN-O5D	-2.58	1.48	1.59
4	C	300	NAP	C4N-C3N	2.50	1.43	1.39
4	C	300	NAP	C3N-C7N	2.49	1.54	1.50
4	C	300	NAP	C2D-C1D	-2.46	1.50	1.53
4	C	300	NAP	C5N-C4N	2.43	1.44	1.38
4	C	300	NAP	O4B-C1B	2.03	1.43	1.41

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	300	NAP	N3A-C2A-N1A	-6.58	118.39	128.68
3	A	400	NAD	O4B-C1B-C2B	-6.34	97.66	106.93
3	A	400	NAD	O5D-C5D-C4D	4.84	125.64	108.99
4	C	300	NAP	C1B-N9A-C4A	-4.45	118.82	126.64
3	A	400	NAD	C5N-C6N-N1N	-4.31	114.21	120.40
3	A	400	NAD	C6N-C5N-C4N	4.29	125.67	119.44
4	C	300	NAP	C3B-C2B-C1B	-4.10	95.18	102.89
3	A	400	NAD	C4N-C3N-C7N	-3.77	110.95	121.04
4	C	300	NAP	O2A-PA-O1A	3.70	130.53	112.24
3	A	400	NAD	C3D-C2D-C1D	3.41	106.11	100.98
4	C	300	NAP	C3N-C2N-N1N	-3.04	117.45	120.43
3	A	400	NAD	C2N-C3N-C7N	3.01	128.21	119.46
4	C	300	NAP	C5N-C4N-C3N	2.85	123.72	120.34
3	A	400	NAD	C5N-C4N-C3N	-2.75	117.09	120.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	300	NAP	C2B-C3B-C4B	2.55	107.52	101.99
4	C	300	NAP	O4B-C4B-C3B	-2.54	100.10	105.11
3	A	400	NAD	C5A-C6A-N6A	2.40	124.00	120.35
3	A	400	NAD	C5D-C4D-C3D	-2.32	106.48	115.18
4	C	300	NAP	O3B-C3B-C4B	-2.30	104.41	111.05
3	A	400	NAD	C2N-C3N-C4N	2.27	120.84	118.26
4	C	300	NAP	O2B-C2B-C1B	2.27	118.28	110.10
4	C	300	NAP	C2D-C3D-C4D	2.26	107.04	102.64
4	C	300	NAP	C6N-C5N-C4N	-2.11	116.37	119.44
4	C	300	NAP	O3D-C3D-C4D	-2.10	104.97	111.05
4	C	300	NAP	C5B-C4B-C3B	-2.08	107.40	115.18
3	A	400	NAD	N3A-C2A-N1A	-2.06	125.46	128.68

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	400	NAD	C5B-O5B-PA-O1A
3	A	400	NAD	C5B-O5B-PA-O2A
4	C	300	NAP	C5B-O5B-PA-O1A
4	C	300	NAP	C5B-O5B-PA-O3
4	C	300	NAP	O4D-C1D-N1N-C2N
4	C	300	NAP	O4D-C1D-N1N-C6N
3	A	400	NAD	O4D-C4D-C5D-O5D
4	C	300	NAP	C1B-C2B-O2B-P2B
3	A	400	NAD	C3D-C4D-C5D-O5D
4	C	300	NAP	C3B-C4B-C5B-O5B
4	C	300	NAP	O4B-C4B-C5B-O5B
3	A	400	NAD	C2N-C3N-C7N-O7N
3	A	400	NAD	C2N-C3N-C7N-N7N
4	C	300	NAP	C4D-C5D-O5D-PN
3	A	400	NAD	C4N-C3N-C7N-O7N
3	A	400	NAD	C3B-C4B-C5B-O5B
3	A	400	NAD	C4N-C3N-C7N-N7N
3	A	400	NAD	O4B-C4B-C5B-O5B
3	A	400	NAD	PN-O3-PA-O1A
4	C	300	NAP	PA-O3-PN-O2N
3	A	400	NAD	C5B-O5B-PA-O3
4	C	300	NAP	O4D-C4D-C5D-O5D
4	C	300	NAP	PA-O3-PN-O1N

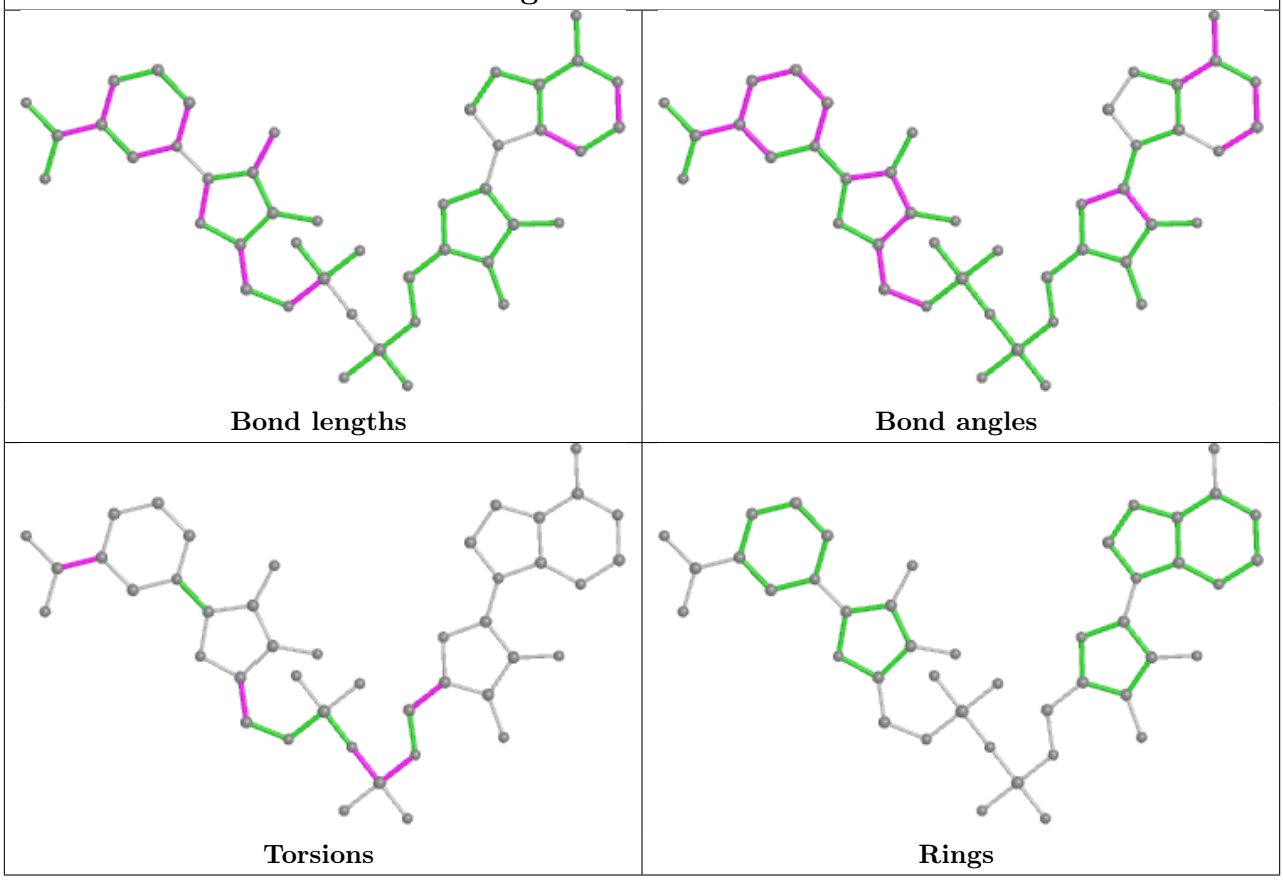
There are no ring outliers.

2 monomers are involved in 6 short contacts:

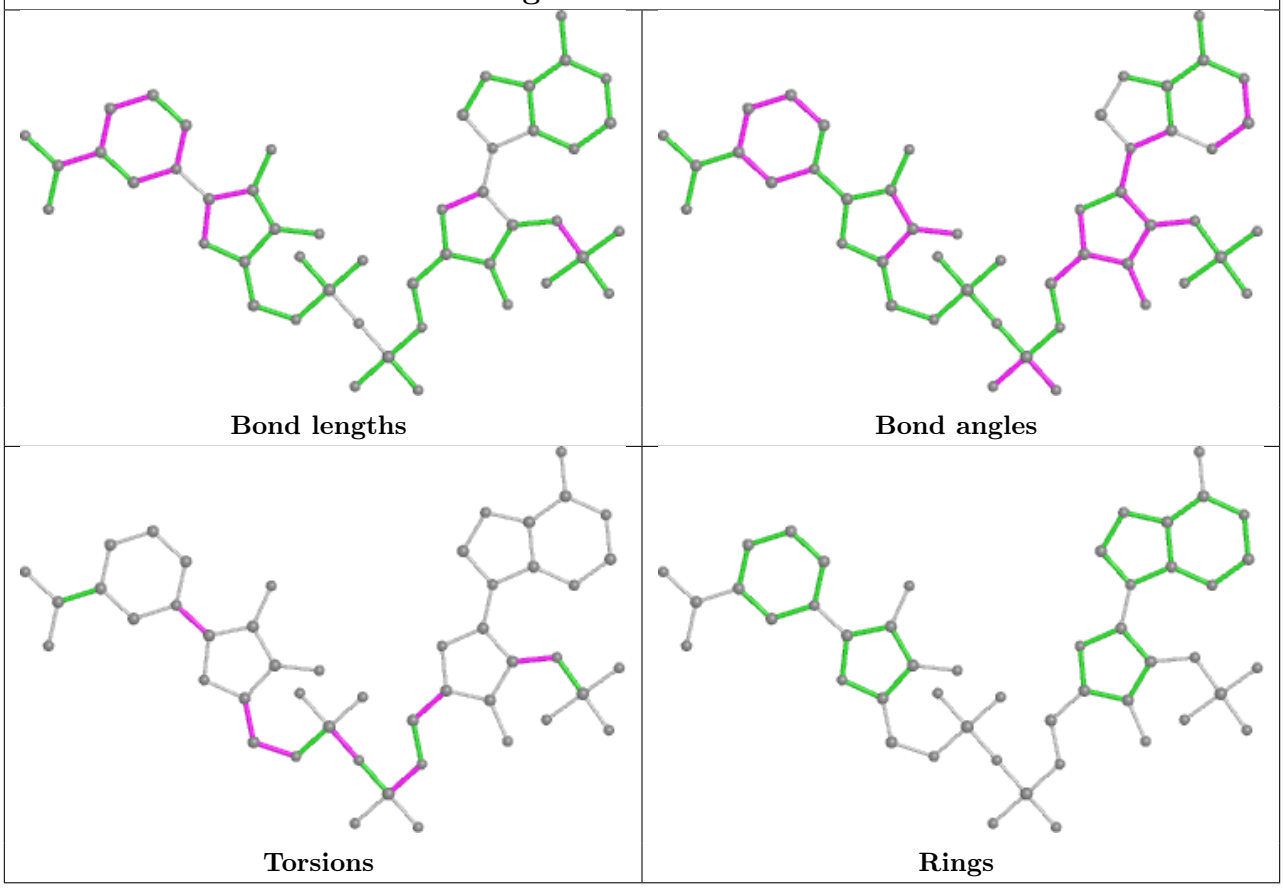
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	NAD	3	0
4	C	300	NAP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand NAD A 400



Ligand NAP C 300



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.