



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:25 am BST

PDB ID : 2BI4  
Title : Lactaldehyde:1,2-propanediol oxidoreductase of Escherichia coli  
Authors : Montella, C.; Bellolell, L.; Badia, J.; Baldoma, L.; Perez, R.; Coll, M.; Aguilar, J.  
Deposited on : 2005-01-20  
Resolution : 2.85 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (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) : 2.11

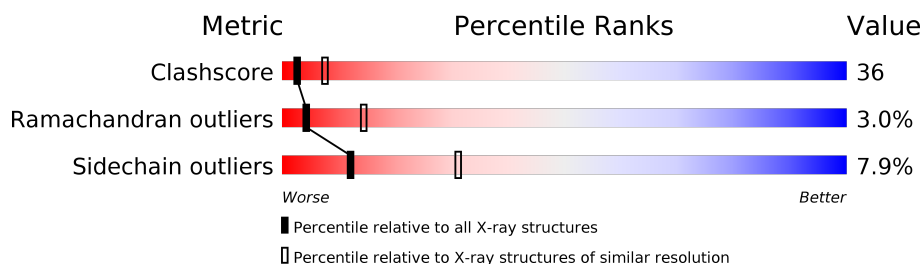
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.85 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	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)

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 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	392	
1	B	392	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5826 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 LACTALDEHYDE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	29	0	0
			2845	1795	493	541	16			
1	B	382	Total	C	N	O	S	38	0	0
			2845	1795	493	541	16			

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Fe	0	0
			1	1		
2	A	1	Total	Fe	0	0
			1	1		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

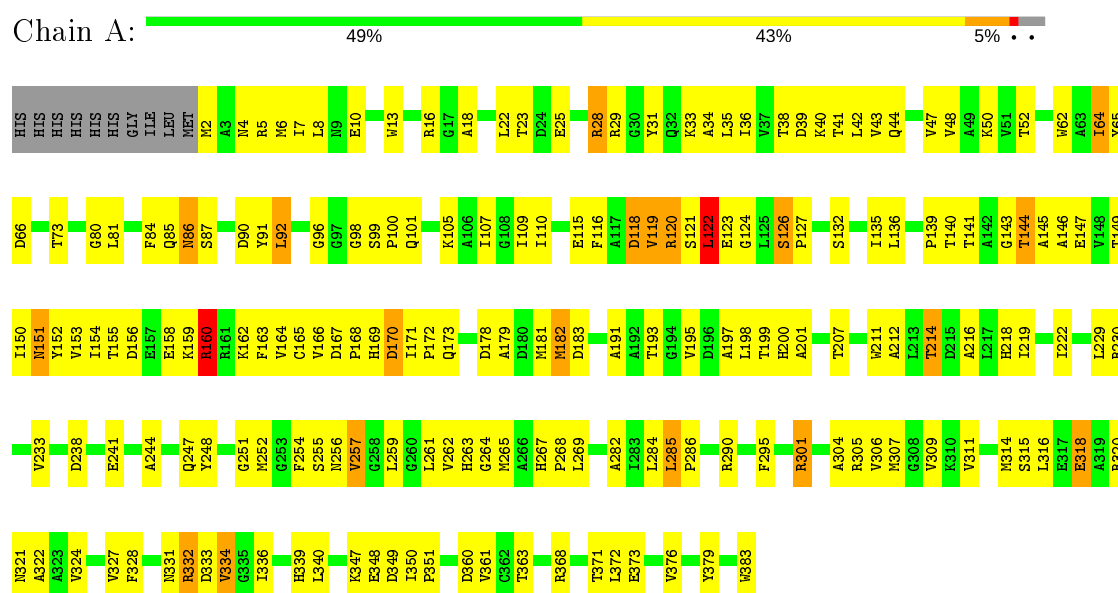
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	20	Total	O	0	0
			20	20		
5	B	25	Total	O	0	0
			25	25		

### 3 Residue-property plots

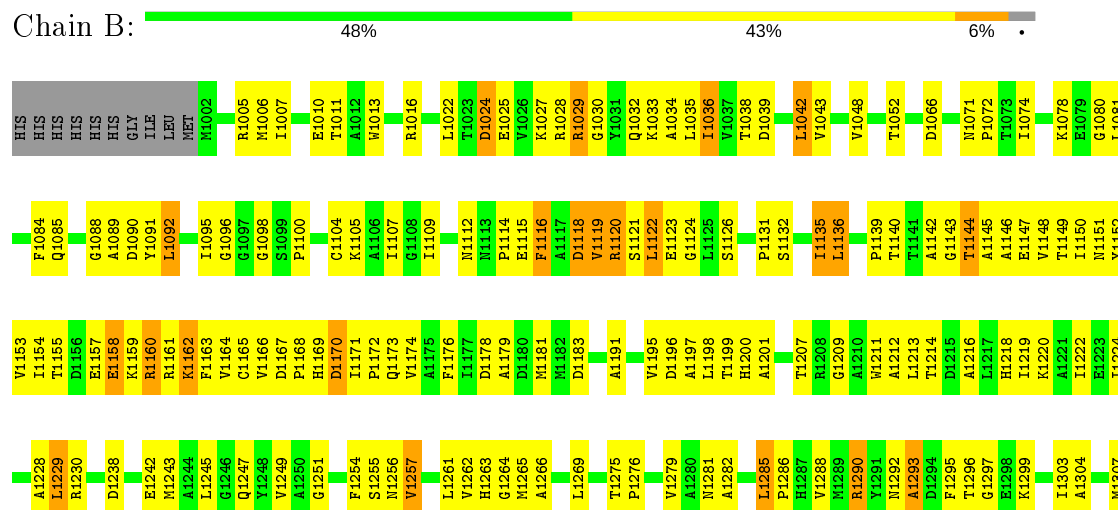
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according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



#### • Molecule 1: LACTALDEHYDE REDUCTASE



G1308	V1309	K1310	V1311	E1312	G1313	M1314	S1315	L1316	E1317	E1318	A1319	R1320	N1321	A1322	A1323	V1324		V1327	F1328	A1329		R1332	D1333	V1334		L1340		I1350	P1351		L1358	D1359	D1360	V1361		R1368		T1371	L1372	E1373	D1374	I1375	V1376	E1377		M1383
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## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.42Å 109.42Å 182.46Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.00 – 2.85	Depositor
% Data completeness (in resolution range)	100.0 (28.00-2.85)	Depositor
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.258 , 0.287	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5826	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	55.0	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: FE, NAD, CL

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	0.42	0/2899	0.65	2/3944 (0.1%)
1	B	0.40	0/2899	0.65	1/3944 (0.0%)
All	All	0.41	0/5798	0.65	3/7888 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ARG	N-CA-C	-5.73	95.52	111.00
1	B	1120	ARG	N-CA-C	-5.57	95.96	111.00
1	A	170	ASP	CB-CA-C	-5.46	99.48	110.40

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	2845	0	2840	218	0
1	B	2845	0	2840	200	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	44	0	26	10	0
3	B	44	0	26	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
5	A	20	0	0	0	0
5	B	25	0	0	1	0
All	All	5826	0	5732	411	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 (411) 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:B:1038:THR:HG22	1:B:1039:ASP:H	1.07	1.16
1:A:38:THR:HG22	1:A:39:ASP:H	1.04	1.09
1:B:1024:ASP:HB3	1:B:1028:ARG:HH12	1.11	1.09
1:B:1317:GLU:HA	1:B:1317:GLU:OE1	1.48	1.08
1:B:1329:ALA:HA	1:B:1332:ARG:NH2	1.67	1.07
1:B:1105:LYS:HD2	1:B:1170:ASP:OD1	1.56	1.05
1:A:38:THR:HG22	1:A:39:ASP:N	1.75	1.01
1:B:1028:ARG:HG3	1:B:1028:ARG:HH11	1.31	0.94
1:A:121:SER:O	1:A:122:LEU:HD12	1.66	0.94
1:A:261:LEU:HD22	1:A:265:MET:HG3	1.50	0.91
1:A:38:THR:CG2	1:A:39:ASP:H	1.81	0.91
1:A:122:LEU:HD23	1:A:152:TYR:OH	1.72	0.90
1:A:191:ALA:HB1	1:A:334:VAL:HG13	1.55	0.88
1:B:1214:THR:HG22	5:B:2011:HOH:O	1.74	0.88
1:A:285:LEU:HB3	1:A:286:PRO:HD3	1.57	0.87
1:A:328:PHE:HA	1:A:332:ARG:NH2	1.90	0.87
1:A:328:PHE:CA	1:A:332:ARG:NH2	2.38	0.87
1:B:1153:VAL:HG12	1:B:1164:VAL:HG22	1.55	0.86
1:B:1038:THR:HG22	1:B:1039:ASP:N	1.92	0.85
1:A:328:PHE:O	1:A:332:ARG:NH2	2.09	0.84
1:B:1329:ALA:HA	1:B:1332:ARG:HH22	1.43	0.83
1:A:328:PHE:HA	1:A:332:ARG:HH22	1.41	0.83
1:A:144:THR:CG2	1:A:146:ALA:HB2	2.10	0.82
1:B:1191:ALA:HB1	1:B:1334:VAL:CG1	2.09	0.82
1:B:1119:VAL:O	1:B:1119:VAL:HG12	1.80	0.81
1:A:144:THR:HG22	1:A:146:ALA:HB2	1.61	0.81
1:A:105:LYS:HD2	1:A:170:ASP:HB3	1.63	0.80
1:B:1024:ASP:HB3	1:B:1028:ARG:NH1	1.94	0.80
1:A:314:MET:HB3	1:A:318:GLU:HB3	1.65	0.79
1:B:1033:LYS:H	1:B:1090:ASP:HB2	1.47	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:VAL:HG12	1:A:164:VAL:HG22	1.64	0.78
1:A:92:LEU:CD2	1:A:92:LEU:N	2.46	0.78
1:A:116:PHE:HB3	1:A:122:LEU:CD1	2.13	0.78
1:B:1178:ASP:OD1	1:B:1181:MET:HG2	1.84	0.78
1:A:118:ASP:OD1	1:A:120:ARG:N	2.15	0.77
1:A:118:ASP:OD1	1:A:119:VAL:N	2.17	0.77
1:A:25:GLU:O	1:A:29:ARG:HG2	1.83	0.77
1:B:1034:ALA:HB2	1:B:1091:TYR:CZ	2.18	0.77
1:A:34:ALA:HB2	1:A:91:TYR:CZ	2.19	0.77
1:B:1028:ARG:NH1	1:B:1028:ARG:HG3	1.99	0.77
1:A:151:ASN:HB3	1:A:166:VAL:HG22	1.67	0.76
1:B:1285:LEU:HB3	1:B:1286:PRO:HD3	1.66	0.76
1:B:1317:GLU:CA	1:B:1317:GLU:OE1	2.30	0.76
1:A:92:LEU:HD22	1:A:92:LEU:N	2.01	0.76
1:A:140:THR:HB	3:A:1385:NAD:H61A	1.51	0.76
1:A:81:LEU:O	1:A:85:GLN:HG3	1.86	0.75
1:A:119:VAL:O	1:A:119:VAL:HG12	1.84	0.75
1:B:1038:THR:CG2	1:B:1039:ASP:H	1.86	0.73
1:A:92:LEU:H	1:A:92:LEU:CD2	2.01	0.72
1:B:1007:ILE:HD13	1:B:1168:PRO:HB3	1.71	0.72
1:B:1151:ASN:HB3	1:B:1166:VAL:HG22	1.71	0.72
1:A:116:PHE:CB	1:A:122:LEU:CD1	2.66	0.72
1:A:7:ILE:CD1	1:A:168:PRO:HB3	2.19	0.72
1:B:1261:LEU:HD22	1:B:1265:MET:HG3	1.70	0.71
1:B:1005:ARG:HD3	1:B:1007:ILE:HG13	1.72	0.71
1:B:1162:LYS:NZ	3:B:2385:NAD:O2D	2.24	0.71
1:B:1372:LEU:O	1:B:1376:VAL:HG23	1.90	0.71
1:A:178:ASP:OD1	1:A:181:MET:HG2	1.90	0.71
1:A:122:LEU:O	1:A:165:CYS:SG	2.49	0.70
1:B:1032:GLN:HB2	1:B:1090:ASP:OD2	1.91	0.70
1:A:339:HIS:HB3	1:A:383:TRP:O	1.91	0.70
1:B:1039:ASP:O	1:B:1043:VAL:HG23	1.91	0.70
1:A:5:ARG:HD3	1:A:7:ILE:HG13	1.72	0.70
1:A:116:PHE:HB3	1:A:122:LEU:HD12	1.73	0.70
1:A:110:ILE:HD11	1:A:119:VAL:HG22	1.74	0.69
1:A:48:VAL:HG11	1:A:64:ILE:HD11	1.75	0.69
1:A:7:ILE:HD13	1:A:168:PRO:HB3	1.72	0.69
1:B:1118:ASP:OD1	1:B:1120:ARG:N	2.26	0.69
1:A:207:THR:HG21	1:A:257:VAL:HG13	1.75	0.69
1:B:1118:ASP:O	1:B:1119:VAL:HB	1.92	0.68
1:A:328:PHE:CA	1:A:332:ARG:HH21	2.06	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:ARG:NH1	1:A:333:ASP:OD2	2.26	0.68
1:A:371:THR:HG22	1:A:373:GLU:OE1	1.94	0.68
1:A:191:ALA:HB1	1:A:334:VAL:CG1	2.24	0.67
1:A:214:THR:HG23	1:A:218:HIS:CE1	2.29	0.67
1:B:1123:GLU:HG2	1:B:1124:GLY:H	1.59	0.67
1:A:328:PHE:C	1:A:332:ARG:NH2	2.47	0.67
1:B:1024:ASP:CB	1:B:1028:ARG:HH12	2.00	0.67
1:B:1038:THR:HG23	1:B:1095:ILE:O	1.94	0.67
1:B:1170:ASP:OD1	1:B:1170:ASP:C	2.33	0.67
1:A:295:PHE:CE1	1:A:368:ARG:HD2	2.29	0.67
1:B:1116:PHE:CB	1:B:1122:LEU:HD13	2.26	0.66
1:A:121:SER:O	1:A:122:LEU:CD1	2.43	0.66
1:A:123:GLU:HG2	1:A:124:GLY:H	1.61	0.66
1:A:286:PRO:HG3	1:A:327:VAL:HG12	1.77	0.65
1:A:290:ARG:HH11	1:A:290:ARG:HG3	1.62	0.65
1:A:120:ARG:NH2	1:A:156:ASP:OD2	2.30	0.65
1:A:164:VAL:HB	1:A:361:VAL:HG21	1.78	0.65
1:A:33:LYS:HD3	1:A:87:SER:O	1.97	0.65
1:A:315:SER:H	1:A:318:GLU:HB2	1.61	0.64
1:B:1196:ASP:OD2	3:B:2385:NAD:H5N	1.97	0.64
1:B:1118:ASP:OD1	1:B:1119:VAL:N	2.31	0.64
1:A:164:VAL:HB	1:A:361:VAL:CG2	2.28	0.64
1:A:13:TRP:CH2	1:B:1005:ARG:HB2	2.32	0.64
1:A:286:PRO:HG3	1:A:327:VAL:CG1	2.27	0.64
1:B:1329:ALA:HA	1:B:1332:ARG:CZ	2.28	0.63
1:A:144:THR:HG22	1:A:146:ALA:CB	2.29	0.63
1:B:1153:VAL:CG1	1:B:1164:VAL:HG22	2.28	0.63
1:A:160:ARG:HG3	1:A:160:ARG:O	1.97	0.63
1:A:6:MET:HA	1:A:256:ASN:OD1	1.99	0.63
1:B:1116:PHE:HB3	1:B:1121:SER:O	1.99	0.62
1:B:1160:ARG:O	1:B:1160:ARG:HG3	1.99	0.62
1:A:372:LEU:O	1:A:376:VAL:HG23	1.98	0.62
1:A:350:ILE:HB	1:A:351:PRO:HD3	1.81	0.62
1:A:118:ASP:O	1:A:119:VAL:HB	1.98	0.62
1:B:1123:GLU:CD	1:B:1161:ARG:HH22	2.03	0.62
1:B:1314:MET:HB3	1:B:1318:GLU:HB3	1.82	0.62
1:A:116:PHE:HB2	1:A:122:LEU:HD11	1.82	0.61
1:B:1092:LEU:N	1:B:1092:LEU:CD2	2.63	0.61
1:B:1119:VAL:CG1	1:B:1119:VAL:O	2.47	0.61
1:A:311:VAL:HA	1:A:314:MET:SD	2.41	0.61
1:A:147:GLU:OE2	1:A:247:GLN:NE2	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:ASN:HD22	1:A:86:ASN:N	1.99	0.61
1:A:301:ARG:NH2	1:A:314:MET:O	2.33	0.61
1:B:1116:PHE:HB2	1:B:1122:LEU:HD13	1.82	0.61
1:B:1123:GLU:OE1	1:B:1161:ARG:NH2	2.32	0.61
1:A:123:GLU:HG2	1:A:124:GLY:N	2.16	0.61
1:B:1105:LYS:CD	1:B:1170:ASP:OD1	2.42	0.61
1:A:347:LYS:HA	1:A:350:ILE:HG12	1.83	0.60
1:A:7:ILE:N	1:A:256:ASN:OD1	2.28	0.60
1:A:116:PHE:CB	1:A:122:LEU:HD11	2.32	0.60
1:A:10:GLU:HB2	1:B:1010:GLU:HB2	1.83	0.60
1:B:1143:GLY:O	1:B:1145:ALA:N	2.35	0.60
1:A:309:VAL:HG11	1:A:322:ALA:HB1	1.84	0.59
1:B:1119:VAL:HG12	1:B:1163:PHE:HZ	1.67	0.59
1:B:1309:VAL:HG11	1:B:1322:ALA:HB1	1.83	0.59
1:B:1105:LYS:HD2	1:B:1170:ASP:CG	2.23	0.59
1:B:1109:ILE:HD13	1:B:1132:SER:HB2	1.83	0.59
1:B:1123:GLU:HG2	1:B:1124:GLY:N	2.18	0.59
1:B:1071:ASN:N	1:B:1072:PRO:HD3	2.18	0.59
1:B:1092:LEU:CD2	1:B:1092:LEU:H	2.14	0.58
1:B:1122:LEU:O	1:B:1165:CYS:SG	2.61	0.58
1:B:1286:PRO:HG3	1:B:1327:VAL:HG12	1.84	0.58
1:A:328:PHE:O	1:A:332:ARG:CZ	2.51	0.58
1:A:290:ARG:HG3	1:A:290:ARG:NH1	2.18	0.58
1:A:5:ARG:HB2	1:B:1013:TRP:CZ3	2.38	0.58
1:A:141:THR:HB	1:A:193:THR:HG21	1.86	0.58
1:B:1164:VAL:HB	1:B:1361:VAL:CG2	2.34	0.58
1:A:152:TYR:HE1	1:A:170:ASP:OD1	1.86	0.58
1:B:1119:VAL:O	1:B:1163:PHE:CZ	2.57	0.58
1:A:199:THR:HG23	1:A:285:LEU:HD13	1.84	0.58
1:A:328:PHE:C	1:A:332:ARG:HH21	2.07	0.58
1:A:36:ILE:HB	1:A:64:ILE:HG12	1.85	0.57
1:A:154:ILE:HB	1:A:163:PHE:CZ	2.39	0.57
1:A:92:LEU:H	1:A:92:LEU:HD23	1.68	0.57
1:A:350:ILE:HD12	1:A:379:TYR:HB3	1.85	0.57
1:B:1207:THR:HG21	1:B:1257:VAL:HG13	1.86	0.57
1:A:331:ASN:HB2	1:A:332:ARG:NH1	2.20	0.57
1:B:1191:ALA:HB1	1:B:1334:VAL:HG12	1.85	0.57
1:B:1039:ASP:O	1:B:1066:ASP:HB2	2.04	0.57
1:A:120:ARG:HH22	1:A:156:ASP:CG	2.09	0.57
1:B:1143:GLY:O	1:B:1247:GLN:OE1	2.22	0.56
1:B:1303:ILE:O	1:B:1307:MET:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1286:PRO:HG3	1:B:1327:VAL:CG1	2.35	0.56
1:A:233:VAL:HG21	1:A:334:VAL:HG22	1.88	0.56
1:A:119:VAL:O	1:A:119:VAL:CG1	2.54	0.56
1:B:1118:ASP:OD1	1:B:1118:ASP:C	2.43	0.56
1:B:1142:ALA:HB1	1:B:1243:MET:HE2	1.88	0.56
1:B:1121:SER:O	1:B:1122:LEU:HB2	2.06	0.56
1:B:1126:SER:HB2	1:B:1167:ASP:HB2	1.88	0.56
1:B:1256:ASN:HD22	1:B:1256:ASN:N	2.03	0.56
1:A:120:ARG:NH2	1:A:156:ASP:CG	2.59	0.56
1:A:348:GLU:HG3	1:A:349:ASP:OD1	2.06	0.55
1:B:1038:THR:CG2	1:B:1039:ASP:N	2.61	0.55
1:B:1116:PHE:HD2	1:B:1121:SER:O	1.89	0.55
1:B:1119:VAL:O	1:B:1163:PHE:CE2	2.59	0.55
1:A:118:ASP:OD1	1:A:120:ARG:HG2	2.06	0.55
1:B:1007:ILE:N	1:B:1256:ASN:OD1	2.29	0.55
1:A:159:LYS:O	1:A:160:ARG:C	2.45	0.55
1:B:1276:PRO:HB2	1:B:1279:VAL:HG22	1.90	0.54
1:B:1006:MET:HA	1:B:1256:ASN:OD1	2.07	0.54
1:B:1098:GLY:CA	1:B:1144:THR:HG21	2.37	0.54
1:B:1214:THR:HG21	1:B:1257:VAL:HG11	1.89	0.54
1:A:99:SER:OG	3:A:1385:NAD:O2A	2.23	0.54
1:B:1034:ALA:HA	1:B:1091:TYR:O	2.08	0.54
1:A:47:VAL:O	1:A:50:LYS:HB2	2.08	0.54
1:B:1007:ILE:CD1	1:B:1168:PRO:HB3	2.38	0.54
1:B:1178:ASP:OD1	1:B:1178:ASP:O	2.26	0.54
1:A:109:ILE:HD13	1:A:132:SER:HB2	1.90	0.53
1:A:328:PHE:CB	1:A:332:ARG:HH21	2.21	0.53
1:B:1329:ALA:CA	1:B:1332:ARG:HH22	2.19	0.53
1:B:1118:ASP:O	1:B:1119:VAL:CB	2.56	0.53
1:B:1096:GLY:HA2	1:B:1140:THR:OG1	2.09	0.53
1:A:153:VAL:CG1	3:A:1385:NAD:H71N	2.22	0.53
1:A:256:ASN:HD22	1:A:256:ASN:N	2.05	0.52
1:B:1164:VAL:HB	1:B:1361:VAL:HG21	1.91	0.52
1:B:1025:GLU:O	1:B:1029:ARG:CD	2.58	0.52
1:B:1295:PHE:CE1	1:B:1368:ARG:HD2	2.45	0.52
1:B:1191:ALA:CB	1:B:1334:VAL:CG1	2.85	0.52
1:A:315:SER:N	1:A:318:GLU:HB2	2.24	0.52
1:A:328:PHE:HB3	1:A:332:ARG:HH21	1.75	0.52
1:A:38:THR:CG2	1:A:39:ASP:N	2.49	0.52
1:B:1150:ILE:C	1:B:1170:ASP:OD2	2.48	0.52
1:B:1315:SER:H	1:B:1318:GLU:HB2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1269:LEU:HD21	1:B:1340:LEU:HD22	1.92	0.51
1:B:1139:PRO:HB3	1:B:1147:GLU:OE1	2.10	0.51
1:B:1374:ASP:O	1:B:1377:GLU:HB2	2.11	0.51
1:A:200:HIS:HE1	3:A:1385:NAD:H5N	1.72	0.51
1:A:80:GLY:HA3	1:A:107:ILE:HD11	1.91	0.51
1:B:1200:HIS:CD2	1:B:1263:HIS:CE1	2.99	0.51
1:A:25:GLU:OE1	1:A:29:ARG:NH2	2.38	0.51
1:B:1028:ARG:CG	1:B:1028:ARG:NH1	2.68	0.51
1:B:1081:LEU:O	1:B:1085:GLN:HG3	2.10	0.51
1:A:152:TYR:CE1	1:A:170:ASP:OD1	2.64	0.51
1:B:1092:LEU:HD23	1:B:1092:LEU:H	1.76	0.51
1:A:153:VAL:CG1	1:A:164:VAL:HG22	2.37	0.50
1:B:1022:LEU:HD11	1:B:1136:LEU:HD23	1.92	0.50
1:B:1092:LEU:N	1:B:1092:LEU:HD22	2.27	0.50
1:B:1161:ARG:CZ	1:B:1163:PHE:HB3	2.42	0.50
1:B:1295:PHE:N	1:B:1295:PHE:CD1	2.79	0.50
1:A:118:ASP:C	1:A:118:ASP:OD1	2.48	0.50
1:A:285:LEU:HB3	1:A:286:PRO:CD	2.36	0.50
1:B:1122:LEU:CD2	1:B:1152:TYR:OH	2.60	0.50
1:B:1005:ARG:HD3	1:B:1007:ILE:CG1	2.40	0.50
1:B:1276:PRO:HD2	1:B:1279:VAL:HG21	1.94	0.50
1:A:200:HIS:NE2	3:A:1385:NAD:H4N	2.26	0.50
1:A:304:ALA:O	1:A:309:VAL:HB	2.11	0.50
1:B:1199:THR:HG23	1:B:1285:LEU:HD13	1.94	0.50
1:B:1304:ALA:O	1:B:1309:VAL:HB	2.10	0.50
1:B:1296:THR:O	1:B:1299:LYS:HB2	2.12	0.50
1:A:149:THR:HG21	1:A:254:PHE:CZ	2.47	0.50
1:A:316:LEU:O	1:A:320:ARG:HG3	2.12	0.49
1:B:1154:ILE:HB	1:B:1163:PHE:CZ	2.47	0.49
1:B:1321:ASN:O	1:B:1324:VAL:HG12	2.12	0.49
1:A:140:THR:HB	3:A:1385:NAD:N6A	2.23	0.49
1:A:31:TYR:HB3	1:A:91:TYR:CD1	2.48	0.49
1:B:1116:PHE:CD1	1:B:1116:PHE:N	2.80	0.49
1:A:139:PRO:HB3	1:A:147:GLU:OE1	2.12	0.49
1:A:16:ARG:HH12	1:A:183:ASP:CG	2.15	0.49
1:A:269:LEU:HD21	1:A:340:LEU:HD22	1.94	0.49
1:A:90:ASP:HB3	1:A:91:TYR:HD1	1.77	0.49
1:A:145:ALA:O	1:A:149:THR:HG23	2.12	0.49
1:B:1149:THR:HG22	1:B:1255:SER:HB2	1.95	0.49
1:A:295:PHE:CD1	1:A:295:PHE:N	2.80	0.49
1:A:171:ILE:O	1:A:171:ILE:HG23	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ARG:NH1	1:A:183:ASP:OD2	2.43	0.49
1:B:1016:ARG:HH12	1:B:1183:ASP:CG	2.16	0.49
1:A:118:ASP:O	1:A:119:VAL:CB	2.61	0.48
1:A:254:PHE:HB2	1:A:259:LEU:HD21	1.94	0.48
1:B:1025:GLU:O	1:B:1029:ARG:HG2	2.13	0.48
1:A:143:GLY:O	1:A:145:ALA:N	2.46	0.48
1:A:200:HIS:CD2	1:A:263:HIS:CE1	3.01	0.48
1:A:251:GLY:HA2	1:A:254:PHE:CE2	2.48	0.48
1:A:98:GLY:HA2	1:A:144:THR:HG21	1.95	0.48
1:A:149:THR:HG22	1:A:255:SER:HB2	1.94	0.48
1:A:35:LEU:HD22	1:A:84:PHE:HA	1.96	0.48
1:A:13:TRP:CZ3	1:B:1005:ARG:HB2	2.49	0.48
1:B:1109:ILE:CD1	1:B:1132:SER:HB2	2.43	0.48
1:B:1122:LEU:HD23	1:B:1152:TYR:OH	2.14	0.48
1:B:1116:PHE:CD2	1:B:1121:SER:O	2.65	0.48
1:B:1191:ALA:CB	1:B:1334:VAL:HG13	2.44	0.48
1:A:99:SER:HB2	1:A:100:PRO:CD	2.44	0.48
1:A:36:ILE:HB	1:A:64:ILE:CG1	2.44	0.48
1:A:64:ILE:HG23	1:A:65:TYR:N	2.29	0.47
1:B:1098:GLY:HA2	1:B:1144:THR:HG21	1.94	0.47
1:B:1230:ARG:NH1	1:B:1333:ASP:OD2	2.47	0.47
1:B:1254:PHE:HA	1:B:1257:VAL:HG12	1.94	0.47
1:B:1316:LEU:HD21	1:B:1320:ARG:NH2	2.29	0.47
1:A:40:LYS:O	1:A:43:VAL:HG22	2.14	0.47
1:A:116:PHE:CD1	1:A:116:PHE:N	2.81	0.47
1:A:264:GLY:HA2	1:A:360:ASP:OD2	2.15	0.47
1:B:1116:PHE:CD2	1:B:1122:LEU:HD12	2.50	0.47
1:A:5:ARG:HB2	1:B:1013:TRP:CE3	2.49	0.47
1:A:101:GLN:HE22	1:A:139:PRO:HB3	1.80	0.47
1:A:321:ASN:O	1:A:324:VAL:HG12	2.14	0.47
1:A:261:LEU:O	1:A:265:MET:N	2.36	0.47
1:A:122:LEU:HA	1:A:126:SER:HB2	1.96	0.47
1:A:5:ARG:HD3	1:A:7:ILE:CG1	2.41	0.47
1:B:1131:PRO:HA	1:B:1169:HIS:HB3	1.96	0.47
1:B:1216:ALA:O	1:B:1219:ILE:HG22	2.15	0.47
1:A:149:THR:HB	1:A:151:ASN:ND2	2.30	0.47
1:B:1266:ALA:CB	1:B:1281:ASN:ND2	2.78	0.47
1:A:144:THR:HG22	1:A:146:ALA:N	2.29	0.46
1:A:118:ASP:CG	1:A:120:ARG:HG2	2.36	0.46
1:A:167:ASP:OD2	1:A:169:HIS:HB2	2.15	0.46
1:A:171:ILE:HD12	1:A:172:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1081:LEU:HA	1:B:1107:ILE:HG23	1.97	0.46
1:B:1220:LYS:O	1:B:1224:ILE:HG13	2.15	0.46
1:B:1036:ILE:CD1	1:B:1036:ILE:N	2.78	0.46
1:A:140:THR:HA	1:A:182:MET:HG2	1.97	0.46
1:A:254:PHE:HB2	1:A:259:LEU:CD2	2.46	0.46
1:B:1090:ASP:HB3	1:B:1091:TYR:HD1	1.79	0.46
1:A:115:GLU:HB2	1:A:116:PHE:CD1	2.50	0.46
1:A:216:ALA:HB2	1:B:1220:LYS:HE2	1.97	0.46
1:B:1371:THR:HG21	1:B:1373:GLU:OE1	2.16	0.46
1:A:13:TRP:HB3	1:A:18:ALA:HB1	1.96	0.46
1:B:1092:LEU:HG	1:B:1104:CYS:SG	2.56	0.46
1:A:261:LEU:O	1:A:262:VAL:C	2.53	0.46
1:A:34:ALA:HB2	1:A:91:TYR:CE1	2.51	0.46
1:B:1171:ILE:HG23	1:B:1171:ILE:O	2.16	0.46
1:A:195:VAL:HG21	1:A:336:ILE:CD1	2.47	0.45
1:A:5:ARG:HB2	1:B:1013:TRP:CH2	2.51	0.45
1:B:1276:PRO:O	1:B:1279:VAL:HG22	2.17	0.45
1:A:216:ALA:O	1:A:219:ILE:HG22	2.17	0.45
1:A:85:GLN:C	1:A:86:ASN:HD22	2.20	0.45
1:A:13:TRP:CZ2	1:B:1005:ARG:HB2	2.52	0.45
1:A:254:PHE:HA	1:A:257:VAL:HG12	1.98	0.45
1:B:1025:GLU:O	1:B:1029:ARG:HD3	2.16	0.45
1:B:1016:ARG:NH1	1:B:1183:ASP:OD2	2.49	0.45
1:B:1011:THR:HG22	1:B:1174:VAL:HG13	1.99	0.45
1:A:195:VAL:HG21	1:A:336:ILE:HD12	1.99	0.44
1:A:191:ALA:CB	1:A:334:VAL:HG13	2.39	0.44
1:B:1144:THR:HG22	1:B:1146:ALA:HB2	1.98	0.44
1:B:1245:LEU:O	1:B:1249:VAL:HG23	2.17	0.44
1:A:31:TYR:HB3	1:A:91:TYR:CE1	2.53	0.44
1:A:86:ASN:N	1:A:86:ASN:ND2	2.66	0.44
1:B:1033:LYS:HB3	1:B:1088:GLY:O	2.17	0.44
1:B:1211:TRP:CH2	1:B:1213:LEU:HB3	2.53	0.44
1:A:178:ASP:O	1:A:178:ASP:OD1	2.35	0.44
1:A:4:ASN:ND2	1:A:211:TRP:CG	2.85	0.44
1:A:201:ALA:HB1	1:A:222:ILE:HG13	1.98	0.44
1:B:1139:PRO:HD2	1:B:1176:PHE:O	2.17	0.44
1:A:144:THR:HG21	1:A:146:ALA:HB2	1.98	0.44
1:B:1048:VAL:HG12	1:B:1052:THR:HG23	1.99	0.44
1:B:1168:PRO:O	1:B:1171:ILE:HG22	2.18	0.44
1:A:126:SER:HA	1:A:127:PRO:HD2	1.78	0.44
1:A:144:THR:HG22	1:A:146:ALA:H	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1074:ILE:HG22	1:B:1078:LYS:HE3	1.99	0.44
1:B:1350:ILE:N	1:B:1351:PRO:HD2	2.32	0.44
1:B:1159:LYS:O	1:B:1160:ARG:C	2.55	0.44
1:B:1161:ARG:O	1:B:1161:ARG:HG3	2.17	0.44
1:A:48:VAL:HG12	1:A:52:THR:HG23	2.00	0.43
1:A:179:ALA:HB2	1:A:241:GLU:HA	2.00	0.43
1:A:8:LEU:HD11	1:A:252:MET:HE2	2.01	0.43
1:A:321:ASN:HA	1:A:324:VAL:HG12	2.01	0.43
1:A:334:VAL:O	1:A:334:VAL:HG13	2.17	0.43
1:A:349:ASP:O	1:A:350:ILE:C	2.56	0.43
1:A:334:VAL:HG12	1:A:336:ILE:HG13	2.00	0.43
1:B:1115:GLU:HB2	1:B:1116:PHE:CD1	2.54	0.43
1:B:1154:ILE:HG22	1:B:1155:THR:N	2.33	0.43
1:B:1104:CYS:SG	1:B:1135:ILE:HD12	2.59	0.43
1:B:1144:THR:O	1:B:1145:ALA:HB3	2.19	0.43
1:A:39:ASP:OD2	3:A:1385:NAD:H2B	2.19	0.43
1:A:140:THR:O	3:A:1385:NAD:N6A	2.51	0.43
1:A:282:ALA:HB1	1:A:331:ASN:ND2	2.34	0.43
1:B:1214:THR:HG23	1:B:1218:HIS:CE1	2.54	0.43
1:A:43:VAL:HG23	1:A:44:GLN:N	2.33	0.43
1:A:96:GLY:HA2	1:A:140:THR:OG1	2.19	0.43
1:A:96:GLY:HA3	1:A:100:PRO:HG3	2.01	0.43
1:B:1131:PRO:HA	1:B:1169:HIS:CD2	2.54	0.43
1:A:247:GLN:HE21	1:A:248:TYR:N	2.16	0.42
1:A:267:HIS:HB2	1:A:268:PRO:HD3	2.01	0.42
1:A:84:PHE:CE1	1:A:92:LEU:HD13	2.54	0.42
1:B:1144:THR:CG2	1:B:1146:ALA:HB2	2.49	0.42
1:B:1251:GLY:HA2	1:B:1254:PHE:CE2	2.54	0.42
1:B:1275:THR:HG22	1:B:1279:VAL:HG23	2.02	0.42
1:A:154:ILE:HG22	1:A:155:THR:N	2.34	0.42
1:B:1039:ASP:OD2	1:B:1042:LEU:HD22	2.19	0.42
1:A:105:LYS:CD	1:A:170:ASP:HB3	2.42	0.42
1:A:256:ASN:N	1:A:256:ASN:ND2	2.68	0.42
1:B:1209:GLY:C	1:B:1257:VAL:CG2	2.88	0.42
1:B:1310:LYS:C	1:B:1312:GLU:H	2.23	0.42
1:A:282:ALA:HB1	1:A:331:ASN:HD22	1.85	0.42
1:B:1292:ASN:O	1:B:1293:ALA:C	2.57	0.42
1:B:1296:THR:O	1:B:1297:GLY:C	2.57	0.42
1:A:149:THR:HA	1:A:255:SER:HB2	2.01	0.42
1:B:1116:PHE:CG	1:B:1122:LEU:CD1	3.02	0.42
1:B:1135:ILE:HG23	1:B:1172:PRO:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1153:VAL:CG1	3:B:2385:NAD:H71N	2.33	0.42
1:B:1201:ALA:HB1	1:B:1222:ILE:HG13	2.02	0.42
1:A:116:PHE:HB2	1:A:122:LEU:CD1	2.44	0.42
1:A:182:MET:CE	1:A:244:ALA:HB2	2.50	0.42
1:B:1027:LYS:C	1:B:1029:ARG:H	2.22	0.42
1:B:1080:GLY:O	1:B:1081:LEU:C	2.57	0.42
1:B:1290:ARG:NH1	1:B:1290:ARG:HG3	2.35	0.42
3:B:2385:NAD:H6N	3:B:2385:NAD:H52N	2.00	0.42
1:B:1149:THR:HG21	1:B:1254:PHE:CZ	2.54	0.42
1:B:1264:GLY:HA2	1:B:1360:ASP:OD2	2.19	0.42
1:A:261:LEU:HG	1:A:363:THR:HG23	2.01	0.41
1:B:1096:GLY:HA3	1:B:1100:PRO:HG3	2.02	0.41
1:B:1290:ARG:HH11	1:B:1290:ARG:HG3	1.85	0.41
1:A:119:VAL:O	1:A:163:PHE:CZ	2.73	0.41
1:A:123:GLU:CG	1:A:124:GLY:H	2.28	0.41
1:A:153:VAL:HG11	3:A:1385:NAD:H71N	1.86	0.41
1:A:73:THR:HA	1:A:155:THR:O	2.20	0.41
1:B:1027:LYS:C	1:B:1029:ARG:N	2.74	0.41
1:B:1218:HIS:NE2	1:B:1254:PHE:N	2.69	0.41
1:B:1358:LEU:HD22	1:B:1372:LEU:HB2	2.02	0.41
1:B:1032:GLN:CB	1:B:1090:ASP:OD2	2.65	0.41
1:A:123:GLU:CG	1:A:124:GLY:N	2.81	0.41
1:A:116:PHE:CE2	1:A:127:PRO:HG2	2.55	0.41
1:A:98:GLY:CA	1:A:144:THR:HG21	2.49	0.41
1:B:1028:ARG:CG	1:B:1028:ARG:HH11	2.09	0.41
1:B:1112:ASN:C	1:B:1114:PRO:HD3	2.39	0.41
1:A:327:VAL:O	1:A:328:PHE:C	2.57	0.41
1:B:1084:PHE:CE1	1:B:1089:ALA:HB3	2.55	0.41
1:B:1148:VAL:O	1:B:1148:VAL:HG22	2.21	0.41
1:A:41:THR:HB	3:A:1385:NAD:O2B	2.21	0.41
1:A:284:LEU:O	1:A:285:LEU:C	2.59	0.41
1:B:1290:ARG:HG2	1:B:1324:VAL:HG21	2.03	0.41
1:B:1035:LEU:CD2	1:B:1084:PHE:HB2	2.51	0.41
1:B:1228:ALA:O	1:B:1229:LEU:C	2.59	0.41
1:B:1262:VAL:HG22	1:B:1288:VAL:HB	2.02	0.41
1:A:40:LYS:HA	1:A:43:VAL:HG22	2.03	0.41
1:A:33:LYS:H	1:A:90:ASP:HB2	1.86	0.41
1:B:1098:GLY:HA3	1:B:1144:THR:HG21	2.02	0.41
1:B:1195:VAL:HG11	1:B:1282:ALA:CB	2.51	0.41
1:A:123:GLU:HG3	1:A:164:VAL:O	2.21	0.41
1:A:38:THR:O	1:A:66:ASP:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1285:LEU:HB3	1:B:1286:PRO:CD	2.44	0.41
1:A:144:THR:CG2	1:A:146:ALA:CB	2.90	0.40
1:A:143:GLY:HA2	1:A:197:ALA:HB2	2.02	0.40
1:A:371:THR:HG22	1:A:372:LEU:N	2.36	0.40
1:A:371:THR:CG2	1:A:373:GLU:OE1	2.65	0.40
1:A:305:ARG:C	1:A:307:MET:H	2.25	0.40
1:B:1238:ASP:O	1:B:1242:GLU:HG3	2.20	0.40
1:B:1256:ASN:ND2	1:B:1256:ASN:N	2.69	0.40
1:A:116:PHE:HE2	1:A:127:PRO:HG2	1.86	0.40
1:A:52:THR:HB	1:A:62:TRP:CZ2	2.56	0.40
1:B:1084:PHE:CE1	1:B:1089:ALA:CB	3.05	0.40
1:B:1143:GLY:HA2	1:B:1197:ALA:HB2	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	380/392 (97%)	330 (87%)	40 (10%)	10 (3%)	5	17
1	B	380/392 (97%)	324 (85%)	43 (11%)	13 (3%)	3	12
All	All	760/784 (97%)	654 (86%)	83 (11%)	23 (3%)	4	14

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	VAL
1	A	122	LEU
1	A	144	THR
1	A	160	ARG
1	A	285	LEU
1	B	1119	VAL

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Mol	Chain	Res	Type
1	B	1122	LEU
1	B	1144	THR
1	B	1160	ARG
1	B	1212	ALA
1	B	1285	LEU
1	B	1030	GLY
1	B	1257	VAL
1	A	212	ALA
1	A	257	VAL
1	B	1158	GLU
1	B	1179	ALA
1	B	1293	ALA
1	A	28	ARG
1	B	1311	VAL
1	B	1312	GLU
1	A	306	VAL
1	A	150	ILE

### 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	291/300 (97%)	264 (91%)	27 (9%)	9	24
1	B	291/300 (97%)	272 (94%)	19 (6%)	17	41
All	All	582/600 (97%)	536 (92%)	46 (8%)	12	31

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	MET
1	A	22	LEU
1	A	23	THR
1	A	28	ARG
1	A	42	LEU
1	A	64	ILE

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Mol	Chain	Res	Type
1	A	86	ASN
1	A	92	LEU
1	A	118	ASP
1	A	122	LEU
1	A	126	SER
1	A	135	ILE
1	A	136	LEU
1	A	151	ASN
1	A	158	GLU
1	A	160	ARG
1	A	162	LYS
1	A	173	GLN
1	A	182	MET
1	A	198	LEU
1	A	214	THR
1	A	229	LEU
1	A	238	ASP
1	A	301	ARG
1	A	318	GLU
1	A	332	ARG
1	A	334	VAL
1	B	1024	ASP
1	B	1029	ARG
1	B	1036	ILE
1	B	1042	LEU
1	B	1092	LEU
1	B	1116	PHE
1	B	1118	ASP
1	B	1135	ILE
1	B	1136	LEU
1	B	1157	GLU
1	B	1158	GLU
1	B	1162	LYS
1	B	1170	ASP
1	B	1173	GLN
1	B	1198	LEU
1	B	1229	LEU
1	B	1290	ARG
1	B	1317	GLU
1	B	1318	GLU

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

Mol	Chain	Res	Type
1	A	32	GLN
1	A	44	GLN
1	A	86	ASN
1	A	101	GLN
1	A	112	ASN
1	A	151	ASN
1	A	281	ASN
1	A	287	HIS
1	B	1071	ASN
1	B	1086	ASN
1	B	1112	ASN
1	B	1173	GLN
1	B	1267	HIS
1	B	1281	ASN
1	B	1331	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAD	A	1385	-	42,48,48	2.11	10 (23%)	50,73,73	1.53	9 (18%)
3	NAD	B	2385	-	42,48,48	2.02	10 (23%)	50,73,73	1.70	13 (26%)

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	1385	-	-	6/26/62/62	0/5/5/5
3	NAD	B	2385	-	-	8/26/62/62	0/5/5/5

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1385	NAD	C2N-N1N	8.08	1.44	1.35
3	B	2385	NAD	C2N-N1N	7.32	1.43	1.35
3	B	2385	NAD	C2A-N3A	4.29	1.39	1.32
3	A	1385	NAD	C4N-C3N	4.23	1.46	1.39
3	B	2385	NAD	C6N-N1N	3.94	1.45	1.35
3	A	1385	NAD	C2A-N1A	3.89	1.41	1.33
3	B	2385	NAD	C2A-N1A	3.73	1.40	1.33
3	A	1385	NAD	C2A-N3A	3.64	1.38	1.32
3	B	2385	NAD	C3N-C7N	3.57	1.55	1.50
3	A	1385	NAD	O4D-C1D	3.20	1.45	1.41
3	A	1385	NAD	C3N-C7N	3.08	1.55	1.50
3	B	2385	NAD	C4N-C3N	3.02	1.44	1.39
3	A	1385	NAD	C6N-N1N	3.00	1.42	1.35
3	A	1385	NAD	C4A-N3A	2.88	1.39	1.35
3	B	2385	NAD	C2N-C3N	2.49	1.42	1.39
3	B	2385	NAD	PA-O1A	-2.46	1.42	1.50
3	A	1385	NAD	C2D-C1D	-2.33	1.50	1.53
3	A	1385	NAD	PA-O1A	-2.23	1.43	1.50
3	B	2385	NAD	O4D-C1D	2.17	1.44	1.41
3	B	2385	NAD	C4A-N3A	2.17	1.38	1.35

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1385	NAD	N3A-C2A-N1A	-4.07	122.32	128.68
3	B	2385	NAD	O4B-C4B-C5B	-4.06	96.02	109.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2385	NAD	N3A-C2A-N1A	-3.99	122.44	128.68
3	B	2385	NAD	C3N-C7N-N7N	3.81	122.32	117.75
3	A	1385	NAD	C3N-C7N-N7N	3.75	122.25	117.75
3	A	1385	NAD	C5A-C6A-N6A	3.58	125.79	120.35
3	A	1385	NAD	O4B-C4B-C5B	-3.51	97.82	109.37
3	B	2385	NAD	C5A-C6A-N6A	3.50	125.67	120.35
3	B	2385	NAD	C2N-C3N-C4N	2.79	121.42	118.26
3	B	2385	NAD	O2A-PA-O1A	2.74	125.81	112.24
3	A	1385	NAD	O2A-PA-O1A	2.52	124.72	112.24
3	B	2385	NAD	O7N-C7N-C3N	-2.51	116.62	119.63
3	B	2385	NAD	O3B-C3B-C4B	-2.43	104.03	111.05
3	A	1385	NAD	O7N-C7N-C3N	-2.35	116.82	119.63
3	A	1385	NAD	C3B-C2B-C1B	2.35	104.51	100.98
3	B	2385	NAD	C2N-N1N-C1D	2.34	124.35	119.14
3	B	2385	NAD	C6N-N1N-C2N	-2.31	119.87	121.97
3	B	2385	NAD	C3B-C2B-C1B	2.27	104.40	100.98
3	B	2385	NAD	C3D-C2D-C1D	2.15	104.21	100.98
3	B	2385	NAD	O4B-C1B-C2B	-2.13	103.81	106.93
3	A	1385	NAD	O3B-C3B-C4B	-2.02	105.21	111.05
3	A	1385	NAD	C1B-N9A-C4A	-2.01	123.11	126.64

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1385	NAD	PA-O3-PN-O5D
3	A	1385	NAD	C5D-O5D-PN-O1N
3	A	1385	NAD	C3D-C4D-C5D-O5D
3	B	2385	NAD	PA-O3-PN-O5D
3	B	2385	NAD	C5D-O5D-PN-O3
3	B	2385	NAD	C5D-O5D-PN-O2N
3	B	2385	NAD	O4D-C4D-C5D-O5D
3	B	2385	NAD	C3D-C4D-C5D-O5D
3	A	1385	NAD	O4D-C4D-C5D-O5D
3	A	1385	NAD	C5D-O5D-PN-O3
3	A	1385	NAD	O4B-C4B-C5B-O5B
3	B	2385	NAD	PA-O3-PN-O1N
3	B	2385	NAD	C5B-O5B-PA-O1A
3	B	2385	NAD	O4B-C4B-C5B-O5B

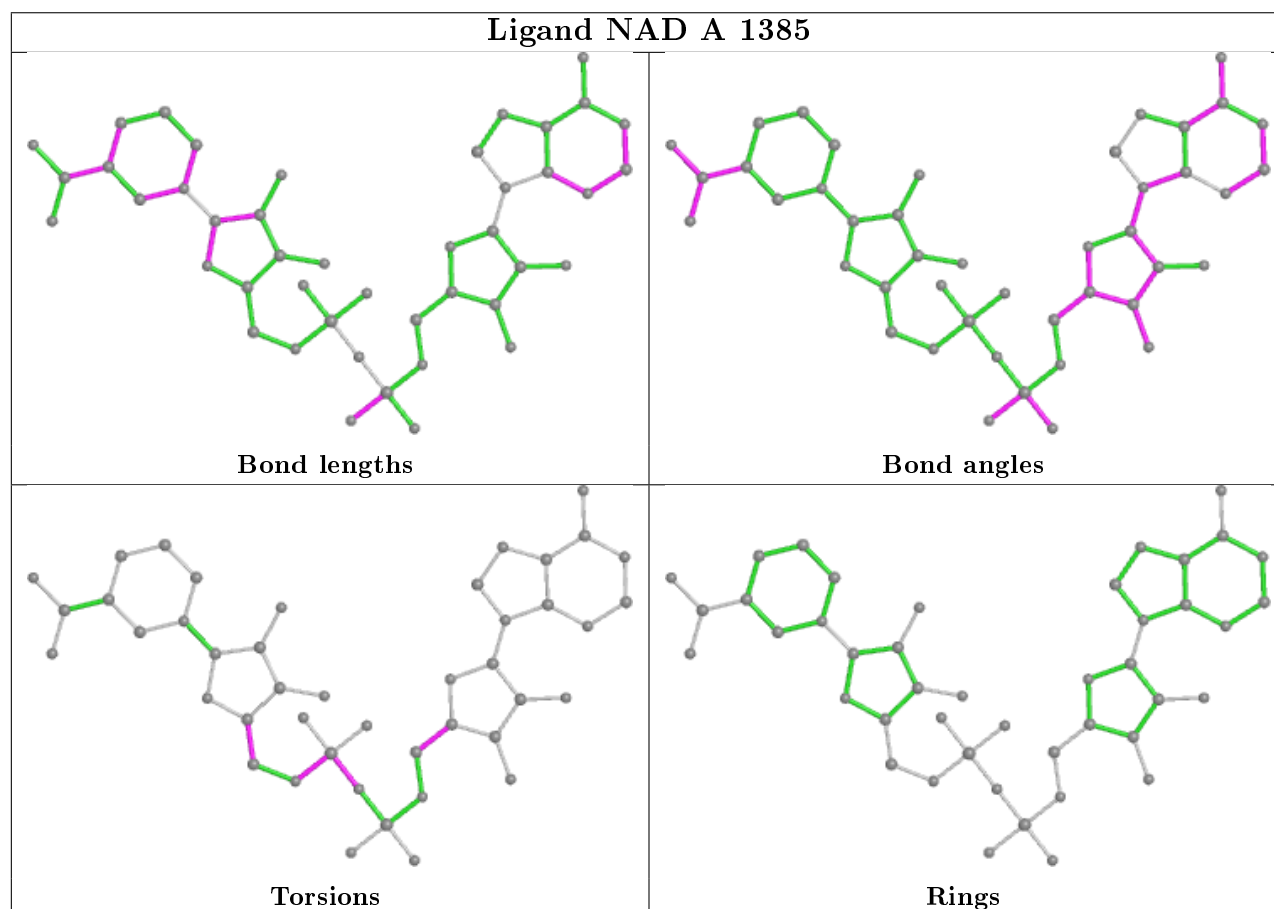
There are no ring outliers.

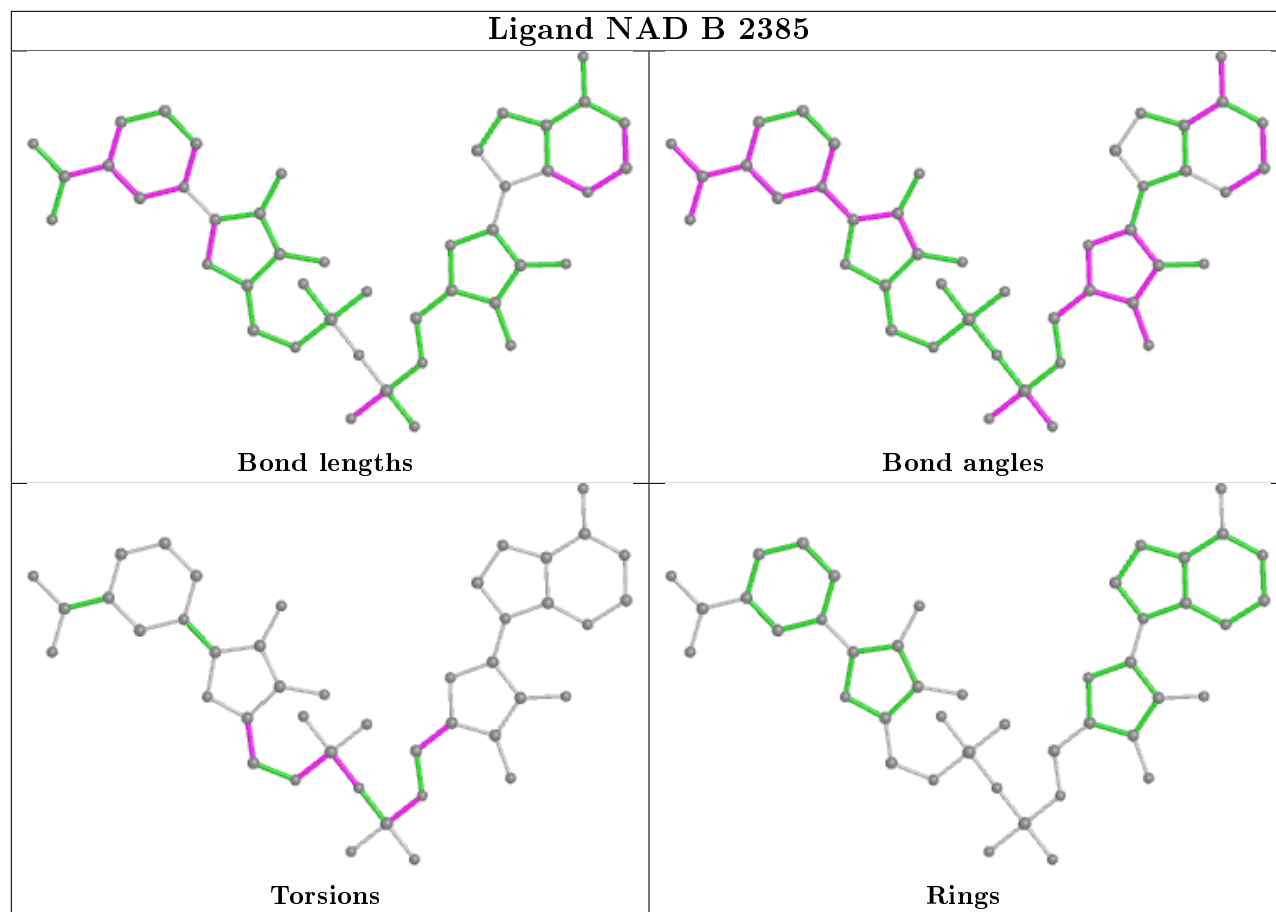
2 monomers are involved in 14 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1385	NAD	10	0
3	B	2385	NAD	4	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 than 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.





## 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

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.