



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:26 PM GMT

PDB ID : 1P1I
Title : Crystal structure of the NAD⁺-bound 1L-myo-inositol 1-phosphate synthase
Authors : Jin, X.; Geiger, J.H.
Deposited on : 2003-04-12
Resolution : 2.40 Å(reported)

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

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

MolProbity : 4.02b-467
Mogul : 1.7 (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

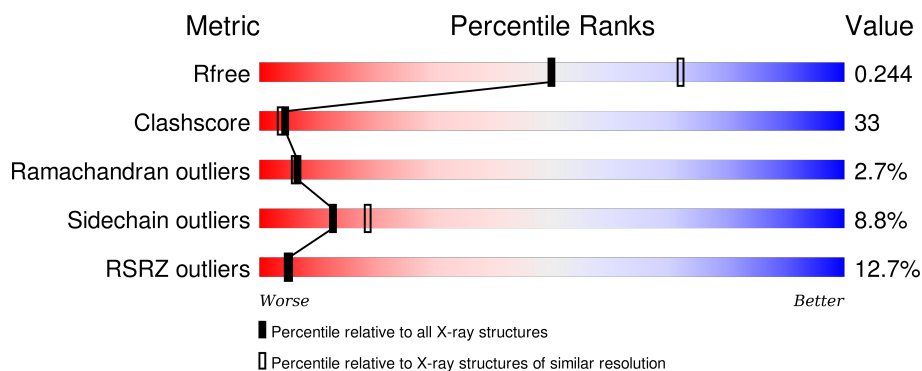
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	533	<div> <div>11%</div> <div>49%</div> <div>41%</div> <div>5%</div> <div>5%</div> </div>
1	B	533	<div> <div>14%</div> <div>43%</div> <div>44%</div> <div>6%</div> <div>7%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8163 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.

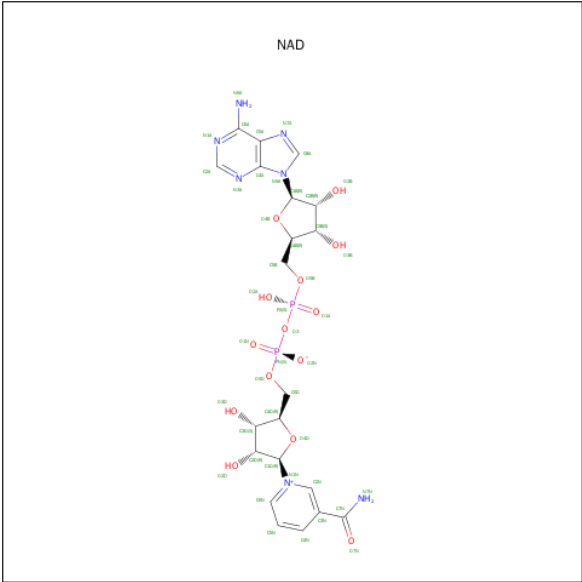
- Molecule 1 is a protein called Inositol-3-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	506	Total	C	N	O	S	0	0	0
			3989	2539	668	766	16			
1	B	498	Total	C	N	O	S	0	0	0
			3926	2502	658	750	16			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ARG	SEE REMARK 999	UNP P11986
A	14	VAL	LEU	SEE REMARK 999	UNP P11986
A	?	-	PHE	SEE REMARK 999	UNP P11986
A	60	LEU	GLU	SEE REMARK 999	UNP P11986
A	?	-	ALA	SEE REMARK 999	UNP P11986
A	98	GLU	LYS	SEE REMARK 999	UNP P11986
A	140	ASN	LYS	SEE REMARK 999	UNP P11986
A	141	ASP	HIS	SEE REMARK 999	UNP P11986
A	201	ASN	GLN	SEE REMARK 999	UNP P11986
A	444	PRO	ALA	SEE REMARK 999	UNP P11986
B	?	-	ARG	SEE REMARK 999	UNP P11986
B	14	VAL	LEU	SEE REMARK 999	UNP P11986
B	?	-	PHE	SEE REMARK 999	UNP P11986
B	60	LEU	GLU	SEE REMARK 999	UNP P11986
B	?	-	ALA	SEE REMARK 999	UNP P11986
B	98	GLU	LYS	SEE REMARK 999	UNP P11986
B	140	ASN	LYS	SEE REMARK 999	UNP P11986
B	141	ASP	HIS	SEE REMARK 999	UNP P11986
B	201	ASN	GLN	SEE REMARK 999	UNP P11986
B	444	PRO	ALA	SEE REMARK 999	UNP P11986

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

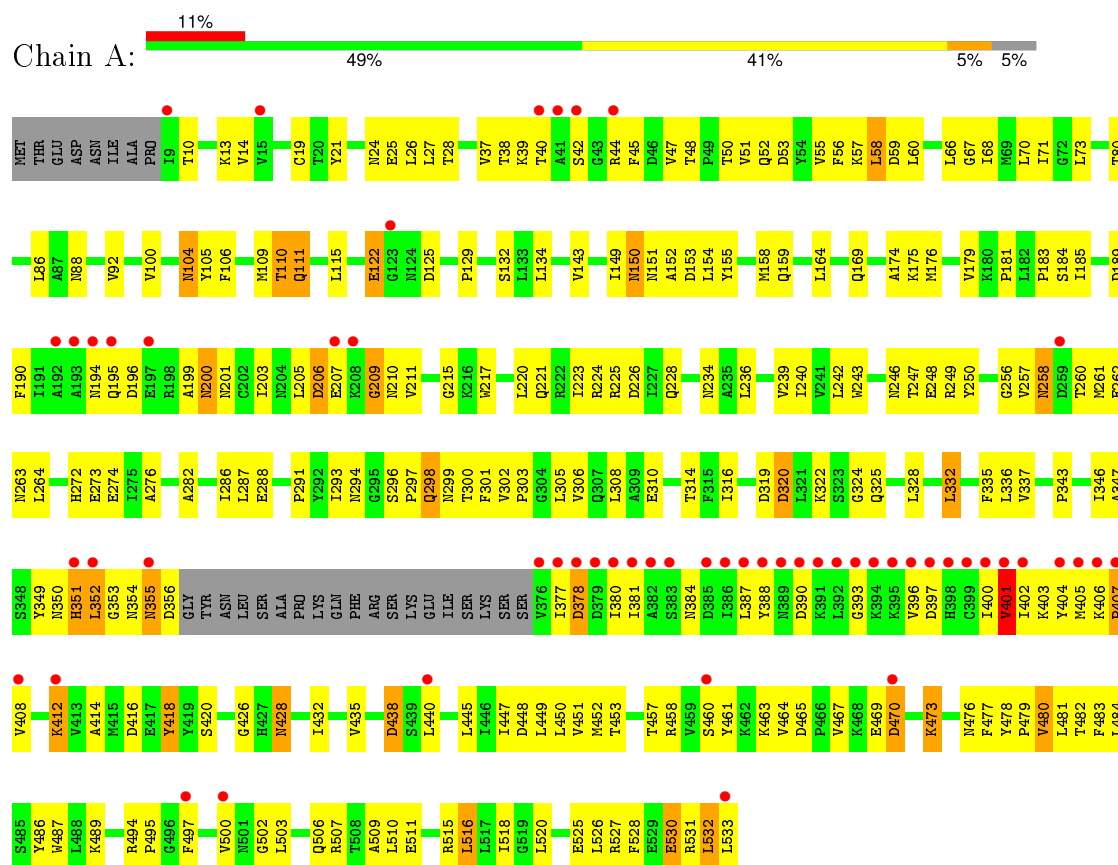
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	90	Total	O	0	0
			90	90		
3	B	70	Total	O	0	0
			70	70		

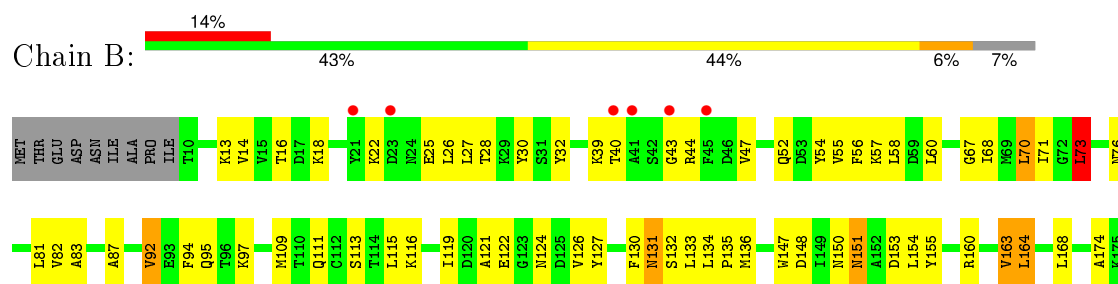
3 Residue-property plots [i](#)

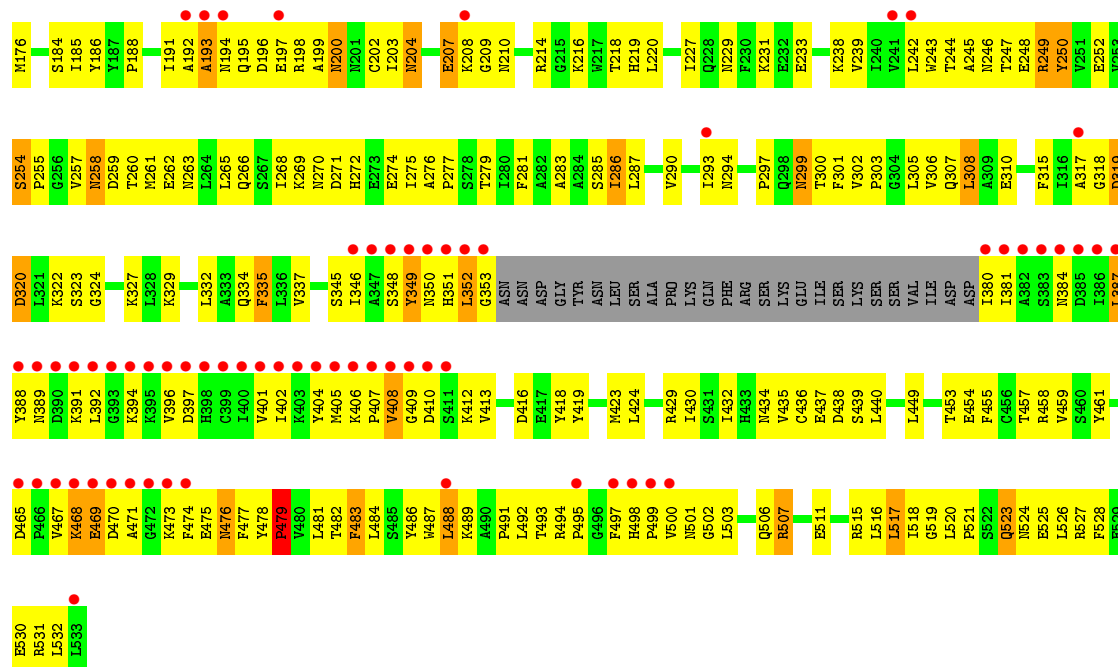
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Inositol-3-phosphate synthase



• Molecule 1: Inositol-3-phosphate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	152.38 Å 97.34 Å 122.89 Å 90.00° 126.53° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 32.92 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.40) 99.4 (32.92-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.39 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.209 , 0.288 0.249 , 0.244	Depositor DCC
R_{free} test set	2807 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	46.4	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.2	EDS
Estimated twinning fraction	0.004 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 56426 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8163	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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	0.39	0/4066	0.65	0/5514
1	B	0.39	0/4003	0.65	2/5427 (0.0%)
All	All	0.39	0/8069	0.65	2/10941 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	73	LEU	CA-CB-CG	5.34	127.58	115.30
1	B	319	ASP	N-CA-C	5.28	125.25	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	3989	0	3994	263	0
1	B	3926	0	3939	303	0
2	A	44	0	26	2	0
2	B	44	0	26	5	0
3	A	90	0	0	6	0
3	B	70	0	0	5	0
All	All	8163	0	7985	525	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 33.

All (525) 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:332:LEU:HD13	1:B:432:ILE:HD11	1.24	1.11
1:A:246:ASN:HD22	2:A:650:NAD:H8A	1.23	1.03
1:B:286:ILE:HD12	1:B:308:LEU:HD13	1.45	0.98
1:B:293:ILE:HA	1:B:317:ALA:HB2	1.49	0.93
1:A:110:THR:HB	1:A:448:ASP:OD1	1.71	0.91
1:B:476:ASN:HD22	1:B:476:ASN:H	1.15	0.90
1:A:104:ASN:ND2	1:A:106:PHE:H	1.69	0.90
1:A:509:ALA:HB1	1:B:531:ARG:HH12	1.38	0.89
1:B:254:SER:H	1:B:258:ASN:HD21	1.22	0.88
1:A:531:ARG:HD3	1:B:482:THR:OG1	1.73	0.87
1:A:353:GLY:HA2	1:A:406:LYS:HA	1.56	0.86
1:A:13:LYS:HE3	1:B:44:ARG:NH1	1.92	0.84
1:A:110:THR:HG22	1:A:111:GLN:NE2	1.93	0.84
1:B:87:ALA:HA	1:B:92:VAL:CG1	2.07	0.84
1:B:468:LYS:HZ2	1:B:471:ALA:HB1	1.41	0.83
1:A:68:ILE:HD12	1:A:450:LEU:HD13	1.61	0.83
1:A:325:GLN:NE2	1:A:350:ASN:HD21	1.76	0.83
1:B:293:ILE:HA	1:B:317:ALA:CB	2.07	0.83
1:A:257:VAL:HG11	1:A:274:GLU:OE2	1.79	0.82
1:B:87:ALA:HA	1:B:92:VAL:HG13	1.62	0.82
1:B:324:GLY:HA2	1:B:327:LYS:HD3	1.62	0.81
1:A:104:ASN:HD22	1:A:106:PHE:H	1.27	0.81
1:A:37:VAL:HG11	1:B:126:VAL:HG11	1.61	0.81
1:B:58:LEU:HD23	1:B:459:VAL:HG22	1.64	0.80
1:A:449:LEU:O	1:A:453:THR:HG23	1.81	0.80
1:B:349:TYR:HB3	1:B:401:VAL:HB	1.64	0.79
1:B:381:ILE:HG23	1:B:388:TYR:HB3	1.65	0.79
1:A:412:LYS:HZ2	1:A:412:LYS:HB3	1.46	0.79
1:B:476:ASN:H	1:B:476:ASN:ND2	1.81	0.77
1:A:502:GLY:O	1:A:506:GLN:HG3	1.84	0.76
1:A:13:LYS:HE3	1:B:44:ARG:HH12	1.48	0.76
1:A:158:MET:HE1	1:A:176:MET:HG3	1.68	0.75
1:B:238:LYS:NZ	1:B:457:THR:HG21	2.01	0.75
1:B:135:PRO:HG2	1:B:458:ARG:HH22	1.51	0.75
1:B:115:LEU:HD22	1:B:511:GLU:HG3	1.67	0.75
1:B:286:ILE:HD11	1:B:308:LEU:HD22	1.68	0.74
1:B:286:ILE:CD1	1:B:308:LEU:HD13	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:LEU:HD12	1:A:354:ASN:OD1	1.88	0.73
1:A:13:LYS:CE	1:B:44:ARG:HH12	2.01	0.73
1:B:109:MET:O	1:B:113:SER:HB3	1.87	0.73
1:B:523:GLN:HA	1:B:523:GLN:HE21	1.53	0.73
1:A:332:LEU:HG	1:A:432:ILE:HD11	1.70	0.73
1:B:52:GLN:HG3	1:B:54:TYR:CE1	2.24	0.73
1:B:286:ILE:HD13	1:B:287:LEU:N	2.04	0.72
1:B:468:LYS:NZ	1:B:471:ALA:HB1	2.05	0.72
1:A:310:GLU:HA	1:A:479:PRO:HG2	1.72	0.71
1:B:492:LEU:HD13	1:B:493:THR:N	2.06	0.71
1:A:352:LEU:H	1:A:352:LEU:HD23	1.53	0.71
1:B:384:ASN:HD21	1:B:387:LEU:HD22	1.55	0.71
1:B:135:PRO:HG2	1:B:458:ARG:NH2	2.04	0.71
1:B:476:ASN:HD22	1:B:476:ASN:N	1.84	0.71
1:B:150:ASN:HD21	1:B:160:ARG:HH12	1.38	0.71
1:A:110:THR:HG22	1:A:111:GLN:HE21	1.55	0.70
1:B:406:LYS:HB3	1:B:407:PRO:HD3	1.73	0.70
1:B:109:MET:HE1	1:B:507:ARG:NH2	2.06	0.70
1:B:410:ASP:OD2	1:B:437:GLU:HB3	1.91	0.70
1:A:104:ASN:HD21	1:B:423:MET:HA	1.56	0.70
1:A:453:THR:O	1:A:457:THR:HG23	1.92	0.70
1:A:533:LEU:HA	1:B:494:ARG:HH22	1.55	0.70
1:B:254:SER:N	1:B:258:ASN:HD21	1.90	0.69
1:A:150:ASN:ND2	1:A:152:ALA:H	1.89	0.69
1:B:261:MET:O	1:B:265:LEU:HD13	1.91	0.69
1:B:302:VAL:O	1:B:306:VAL:HG23	1.93	0.69
1:A:44:ARG:HH21	1:B:13:LYS:HG2	1.58	0.68
1:A:224:ARG:O	1:A:228:GLN:HG3	1.93	0.68
1:B:294:ASN:H	1:B:317:ALA:HB3	1.58	0.68
1:A:353:GLY:CA	1:A:406:LYS:HA	2.23	0.68
1:B:454:GLU:O	1:B:457:THR:HB	1.93	0.67
1:B:131:ASN:HD22	1:B:132:SER:N	1.93	0.67
1:B:345:SER:HB3	1:B:419:TYR:HB3	1.76	0.67
1:B:25:GLU:OE1	1:B:57:LYS:HD3	1.93	0.67
1:A:400:ILE:HG22	1:A:401:VAL:H	1.58	0.67
1:A:272:HIS:CD2	1:A:274:GLU:HB2	2.29	0.67
1:A:473:LYS:H	1:A:473:LYS:HD2	1.60	0.67
1:B:109:MET:HE1	1:B:507:ARG:HG3	1.77	0.67
1:A:352:LEU:N	1:A:352:LEU:HD23	2.09	0.66
1:B:473:LYS:HD2	1:B:474:PHE:H	1.60	0.66
1:B:516:LEU:HD12	1:B:517:LEU:N	2.10	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:384:ASN:ND2	1:B:387:LEU:HD22	2.10	0.66
1:A:328:LEU:HD21	1:B:430:ILE:HG21	1.78	0.66
1:A:246:ASN:ND2	2:A:650:NAD:H8A	2.05	0.66
1:B:507:ARG:HG3	1:B:507:ARG:HH21	1.60	0.65
1:B:227:ILE:O	1:B:231:LYS:HG3	1.96	0.65
1:B:58:LEU:HD23	1:B:459:VAL:CG2	2.24	0.65
1:B:492:LEU:HD13	1:B:493:THR:H	1.61	0.65
1:B:258:ASN:HD22	1:B:258:ASN:H	1.44	0.65
1:B:404:TYR:HE1	1:B:406:LYS:HB2	1.62	0.65
1:B:334:GLN:HA	1:B:380:ILE:HD13	1.78	0.65
1:B:57:LYS:HE3	1:B:474:PHE:CD2	2.33	0.64
1:A:200:ASN:H	1:A:200:ASN:HD22	1.45	0.64
1:A:478:TYR:CE1	1:A:494:ARG:HB3	2.32	0.64
1:A:350:ASN:HB2	1:A:402:ILE:HA	1.78	0.64
1:B:203:ILE:HG13	1:B:204:ASN:N	2.13	0.64
1:A:481:LEU:HB3	1:A:484:LEU:CD2	2.28	0.64
1:B:478:TYR:CD1	1:B:479:PRO:HD2	2.33	0.64
1:A:478:TYR:CD2	1:A:479:PRO:HD2	2.32	0.64
1:A:37:VAL:CG1	1:B:126:VAL:HG11	2.27	0.63
1:B:247:THR:CG2	1:B:297:PRO:HG2	2.28	0.63
1:A:104:ASN:HD22	1:A:105:TYR:N	1.96	0.63
1:B:121:ALA:HB3	1:B:122:GLU:OE2	1.97	0.63
1:A:190:PHE:O	1:A:248:GLU:HA	1.98	0.63
1:B:329:LYS:HG3	1:B:418:TYR:OH	1.98	0.63
1:A:195:GLN:HE22	1:A:355:ASN:HD21	1.48	0.62
1:B:387:LEU:HA	1:B:392:LEU:HD22	1.80	0.62
1:B:277:PRO:HG2	3:B:665:HOH:O	1.99	0.62
1:B:249:ARG:HH21	1:B:249:ARG:HG2	1.65	0.62
1:B:216:LYS:HA	1:B:219:HIS:CD2	2.35	0.62
1:A:481:LEU:HB3	1:A:484:LEU:HD23	1.81	0.62
1:B:299:ASN:HD22	1:B:299:ASN:N	1.97	0.62
1:B:493:THR:HG21	1:B:499:PRO:HD3	1.82	0.62
1:A:149:ILE:O	1:A:199:ALA:HA	1.99	0.62
1:B:40:THR:HG23	1:B:44:ARG:O	1.99	0.61
1:B:30:TYR:CE1	1:B:32:TYR:HB2	2.35	0.61
1:B:352:LEU:HD23	1:B:352:LEU:N	2.15	0.61
1:B:261:MET:H	1:B:307:GLN:NE2	1.99	0.61
1:B:515:ARG:NH1	1:B:521:PRO:O	2.33	0.61
1:B:57:LYS:HB2	1:B:474:PHE:CE2	2.35	0.61
1:A:14:VAL:HA	1:B:47:VAL:HG13	1.81	0.61
1:A:527:ARG:HB2	1:A:531:ARG:NH2	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:ILE:O	1:A:401:VAL:HG23	2.00	0.61
1:B:302:VAL:HB	1:B:303:PRO:HD2	1.83	0.61
1:B:56:PHE:CD2	1:B:461:TYR:HB3	2.34	0.61
1:A:21:TYR:CE1	1:A:26:LEU:HD13	2.36	0.61
1:B:14:VAL:HG11	1:B:518:ILE:O	2.01	0.61
1:B:151:ASN:H	1:B:200:ASN:HD21	1.47	0.61
1:B:315:PHE:CD1	1:B:481:LEU:HD11	2.36	0.60
1:A:486:TYR:HA	1:A:506:GLN:OE1	2.01	0.60
1:A:381:ILE:HG13	1:A:388:TYR:CG	2.36	0.60
1:A:150:ASN:HD22	1:A:152:ALA:H	1.48	0.60
1:A:404:TYR:CE2	1:A:406:LYS:HB2	2.37	0.60
1:B:185:ILE:HG12	1:B:203:ILE:HD11	1.84	0.60
1:B:58:LEU:CD1	1:B:60:LEU:HD23	2.31	0.60
1:A:206:ASP:OD1	1:A:210:ASN:HB2	2.02	0.60
1:A:272:HIS:CD2	1:A:274:GLU:H	2.20	0.60
1:A:217:TRP:O	1:A:221:GLN:HG2	2.01	0.60
1:B:381:ILE:HG23	1:B:388:TYR:CB	2.32	0.59
1:A:405:MET:O	1:A:408:VAL:HG22	2.01	0.59
1:B:469:GLU:HG2	1:B:470:ASP:N	2.17	0.59
1:A:19:CYS:HB3	1:A:28:THR:OG1	2.03	0.59
1:A:316:ILE:HD11	1:A:480:VAL:HG22	1.82	0.59
1:A:500:VAL:HG21	1:B:527:ARG:CZ	2.33	0.59
1:A:509:ALA:CB	1:B:531:ARG:HH12	2.13	0.59
1:A:350:ASN:O	1:A:351:HIS:HB3	2.02	0.59
1:B:349:TYR:CB	1:B:401:VAL:HB	2.31	0.59
1:A:158:MET:CE	1:A:176:MET:HG3	2.32	0.59
1:A:38:THR:HG22	1:A:39:LYS:N	2.18	0.59
1:B:352:LEU:HD21	1:B:402:ILE:HD11	1.86	0.58
1:A:294:ASN:ND2	1:A:296:SER:H	2.01	0.58
1:B:299:ASN:HD22	1:B:299:ASN:H	1.52	0.58
1:A:14:VAL:HG21	1:A:518:ILE:HD12	1.85	0.58
1:A:516:LEU:C	1:A:516:LEU:HD23	2.24	0.58
1:B:200:ASN:ND2	1:B:200:ASN:C	2.58	0.57
1:A:185:ILE:HG12	1:A:203:ILE:HD11	1.86	0.57
1:A:205:LEU:HD13	1:A:209:GLY:HA2	1.85	0.57
1:A:92:VAL:O	1:A:92:VAL:HG12	2.04	0.57
1:B:134:LEU:HD21	1:B:518:ILE:CG2	2.34	0.57
1:A:297:PRO:HB3	1:A:320:ASP:OD2	2.04	0.57
1:B:255:PRO:HA	1:B:259:ASP:CG	2.25	0.57
1:B:58:LEU:HD12	1:B:60:LEU:HD23	1.86	0.57
1:A:44:ARG:NH2	1:B:13:LYS:HG2	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:LEU:HD13	1:B:432:ILE:CD1	2.16	0.57
1:A:533:LEU:HD13	1:B:478:TYR:HD1	1.69	0.57
1:B:249:ARG:HD2	1:B:250:TYR:O	2.05	0.57
1:B:348:SER:HB3	1:B:416:ASP:OD1	2.04	0.57
1:B:192:ALA:O	1:B:194:ASN:N	2.38	0.56
1:B:489:LYS:O	1:B:491:PRO:HD3	2.05	0.56
1:B:26:LEU:O	1:B:57:LYS:HA	2.04	0.56
1:A:125:ASP:CG	1:A:527:ARG:HH22	2.07	0.56
1:B:134:LEU:HD12	1:B:455:PHE:HZ	1.70	0.56
1:B:97:LYS:HB2	1:B:97:LYS:NZ	2.20	0.56
1:A:388:TYR:HA	1:A:393:GLY:O	2.05	0.56
1:B:73:LEU:HD13	1:B:154:LEU:HD11	1.85	0.56
1:A:52:GLN:NE2	1:A:463:LYS:HD2	2.21	0.56
1:B:109:MET:HE1	1:B:507:ARG:HH21	1.69	0.56
1:B:528:PHE:O	1:B:532:LEU:HB2	2.04	0.56
1:A:406:LYS:O	1:A:408:VAL:N	2.34	0.56
1:B:131:ASN:HD22	1:B:132:SER:H	1.52	0.56
1:B:310:GLU:HG2	1:B:479:PRO:HG2	1.87	0.56
1:B:30:TYR:HE1	1:B:32:TYR:HB2	1.71	0.56
1:A:494:ARG:HB2	1:A:495:PRO:HD2	1.88	0.55
1:B:476:ASN:N	1:B:476:ASN:ND2	2.47	0.55
1:A:37:VAL:HG22	1:A:47:VAL:HG22	1.86	0.55
1:B:192:ALA:O	1:B:195:GLN:N	2.40	0.55
1:B:238:LYS:HZ1	1:B:457:THR:HG21	1.71	0.55
1:A:200:ASN:N	1:A:200:ASN:HD22	2.02	0.55
1:A:406:LYS:HD3	1:A:406:LYS:C	2.25	0.55
1:B:323:SER:HB3	1:B:489:LYS:HZ1	1.71	0.55
1:A:39:LYS:HA	1:A:44:ARG:O	2.05	0.55
1:B:220:LEU:C	1:B:220:LEU:HD23	2.27	0.55
1:B:500:VAL:HG12	1:B:502:GLY:H	1.70	0.55
1:A:13:LYS:NZ	1:B:44:ARG:HH12	2.04	0.55
1:B:468:LYS:NZ	1:B:471:ALA:CB	2.69	0.55
1:B:391:LYS:HB2	1:B:391:LYS:HZ2	1.71	0.55
1:A:381:ILE:HD12	1:A:396:VAL:HG23	1.89	0.55
1:A:325:GLN:HE22	1:A:350:ASN:HD21	1.50	0.54
1:A:272:HIS:HD2	1:A:274:GLU:HB2	1.72	0.54
1:B:404:TYR:CE1	1:B:406:LYS:HB2	2.42	0.54
1:A:205:LEU:HB3	1:A:209:GLY:HA2	1.89	0.54
1:B:332:LEU:C	1:B:332:LEU:HD23	2.28	0.54
1:A:377:ILE:O	1:A:381:ILE:HG22	2.07	0.54
1:B:247:THR:HG21	1:B:297:PRO:HG2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:VAL:O	1:A:306:VAL:HG23	2.07	0.54
1:A:412:LYS:O	1:A:435:VAL:HA	2.07	0.54
1:A:528:PHE:CD1	1:A:532:LEU:HD22	2.43	0.54
1:B:247:THR:HG23	1:B:297:PRO:HG2	1.90	0.54
1:B:150:ASN:ND2	1:B:160:ARG:HH12	2.03	0.54
1:B:320:ASP:OD2	2:B:660:NAD:H6N	2.08	0.54
1:A:68:ILE:HG22	1:A:70:LEU:HD13	1.90	0.54
1:B:200:ASN:HD22	1:B:200:ASN:C	2.09	0.54
1:B:255:PRO:HA	1:B:259:ASP:OD1	2.08	0.54
1:A:39:LYS:HD2	1:A:44:ARG:O	2.08	0.54
1:A:412:LYS:HZ1	1:A:414:ALA:HB2	1.74	0.53
1:B:95:GLN:NE2	1:B:95:GLN:HA	2.23	0.53
1:B:272:HIS:CE1	1:B:274:GLU:HB2	2.43	0.53
1:A:402:ILE:HD13	3:A:737:HOH:O	2.07	0.53
1:B:516:LEU:HD12	1:B:516:LEU:C	2.29	0.53
1:A:406:LYS:HD3	1:A:406:LYS:O	2.08	0.53
1:B:468:LYS:HG2	1:B:471:ALA:HB2	1.90	0.53
1:A:350:ASN:N	1:A:401:VAL:O	2.42	0.53
1:A:322:LYS:HD3	1:A:503:LEU:HD13	1.89	0.53
1:A:287:LEU:HD21	1:A:308:LEU:HD11	1.91	0.53
1:B:16:THR:HG21	1:B:519:GLY:HA3	1.91	0.53
1:B:283:ALA:HA	1:B:286:ILE:HG23	1.91	0.52
1:A:58:LEU:HD13	1:A:60:LEU:HD12	1.91	0.52
1:B:244:THR:O	1:B:244:THR:HG23	2.08	0.52
1:A:353:GLY:O	1:A:406:LYS:HG2	2.08	0.52
1:A:282:ALA:O	1:A:286:ILE:HG13	2.08	0.52
1:A:150:ASN:HD22	1:A:151:ASN:N	2.06	0.52
1:A:509:ALA:HB1	1:B:531:ARG:NH1	2.17	0.52
1:A:38:THR:HG22	1:A:39:LYS:H	1.74	0.52
1:A:150:ASN:HA	1:A:200:ASN:ND2	2.24	0.52
1:A:190:PHE:CE2	1:A:276:ALA:HB2	2.44	0.52
1:A:14:VAL:CG2	1:A:518:ILE:HD12	2.39	0.52
1:A:258:ASN:ND2	1:A:258:ASN:H	2.08	0.52
1:B:208:LYS:C	1:B:210:ASN:H	2.13	0.52
1:A:378:ASP:HA	1:A:381:ILE:HG22	1.91	0.52
1:B:238:LYS:HZ2	1:B:457:THR:HG21	1.75	0.51
1:A:40:THR:HB	1:A:44:ARG:HB3	1.91	0.51
1:A:473:LYS:H	1:A:473:LYS:CD	2.23	0.51
1:B:484:LEU:O	1:B:488:LEU:HD22	2.10	0.51
1:B:381:ILE:HD11	1:B:396:VAL:HG23	1.92	0.51
1:A:418:TYR:OH	1:A:432:ILE:HD12	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:389:ASN:ND2	1:B:391:LYS:HB3	2.25	0.51
1:B:299:ASN:ND2	1:B:299:ASN:H	2.07	0.51
1:A:164:LEU:N	1:A:164:LEU:HD12	2.25	0.51
1:A:105:TYR:HA	1:B:423:MET:HE3	1.93	0.51
1:A:528:PHE:C	1:A:530:GLU:H	2.14	0.51
1:B:322:LYS:HA	1:B:489:LYS:HG3	1.93	0.51
1:B:134:LEU:HD21	1:B:518:ILE:HG22	1.93	0.51
1:B:207:GLU:C	1:B:209:GLY:H	2.13	0.50
1:B:28:THR:O	1:B:55:VAL:HA	2.11	0.50
1:A:273:GLU:HB2	3:A:722:HOH:O	2.11	0.50
1:B:523:GLN:CA	1:B:523:GLN:HE21	2.18	0.50
1:B:467:VAL:O	1:B:468:LYS:HB2	2.12	0.50
1:B:306:VAL:O	1:B:310:GLU:HG3	2.11	0.50
1:A:150:ASN:HA	1:A:200:ASN:HD21	1.75	0.50
1:A:484:LEU:HD22	1:A:484:LEU:N	2.27	0.50
1:B:67:GLY:O	1:B:68:ILE:HD13	2.11	0.50
1:B:134:LEU:HB3	1:B:135:PRO:HD2	1.93	0.50
1:A:445:LEU:HD21	1:A:487:TRP:HB3	1.94	0.50
1:A:256:GLY:C	1:A:263:ASN:HD22	2.15	0.50
1:A:469:GLU:HG2	1:A:470:ASP:N	2.26	0.50
1:A:258:ASN:HD22	1:A:258:ASN:H	1.58	0.50
1:B:122:GLU:H	1:B:122:GLU:CD	2.15	0.50
1:B:97:LYS:CB	1:B:97:LYS:NZ	2.75	0.50
1:B:95:GLN:CA	1:B:95:GLN:HE21	2.25	0.50
1:A:88:ASN:HD21	1:A:104:ASN:HA	1.76	0.49
1:A:150:ASN:C	1:A:150:ASN:HD22	2.15	0.49
1:B:285:SER:O	1:B:290:VAL:HG22	2.12	0.49
1:A:100:VAL:HG13	1:A:100:VAL:O	2.11	0.49
1:A:426:GLY:HA3	1:B:440:LEU:CD1	2.42	0.49
1:B:332:LEU:CD1	1:B:432:ILE:HD11	2.17	0.49
1:A:435:VAL:HG13	1:B:429:ARG:HB3	1.94	0.49
1:B:389:ASN:HD21	1:B:391:LYS:HB3	1.77	0.49
1:B:258:ASN:HD22	1:B:258:ASN:N	2.03	0.49
1:A:426:GLY:HA3	1:B:440:LEU:HD13	1.94	0.49
1:B:413:VAL:HA	1:B:434:ASN:O	2.12	0.49
1:B:337:VAL:HG21	1:B:380:ILE:HG22	1.95	0.49
1:A:264:LEU:HD22	1:A:305:LEU:HA	1.93	0.49
1:A:384:ASN:CG	1:A:387:LEU:HD12	2.33	0.49
1:A:104:ASN:HD22	1:A:106:PHE:N	2.03	0.49
1:A:478:TYR:HE1	1:A:494:ARG:HB3	1.77	0.49
1:B:350:ASN:HD21	1:B:402:ILE:HG13	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:405:MET:O	1:B:408:VAL:HG23	2.12	0.49
1:B:381:ILE:HD11	1:B:396:VAL:CG2	2.43	0.49
1:B:252:GLU:O	1:B:274:GLU:OE1	2.29	0.49
1:B:337:VAL:HG21	1:B:380:ILE:CG2	2.43	0.49
1:A:347:ALA:HB1	1:A:349:TYR:CE2	2.48	0.49
1:B:348:SER:CB	1:B:416:ASP:HA	2.43	0.49
1:A:351:HIS:HA	1:A:403:LYS:O	2.12	0.48
1:A:335:PHE:HZ	1:B:327:LYS:HE2	1.77	0.48
1:B:192:ALA:C	1:B:194:ASN:H	2.16	0.48
1:A:343:PRO:HA	1:A:420:SER:HA	1.94	0.48
1:B:249:ARG:NH2	1:B:249:ARG:HG2	2.27	0.48
1:B:218:THR:HG23	3:B:699:HOH:O	2.12	0.48
1:A:24:ASN:O	1:A:60:LEU:HB2	2.12	0.48
1:B:219:HIS:HE1	3:B:697:HOH:O	1.95	0.48
1:B:449:LEU:HD21	1:B:487:TRP:HB2	1.96	0.48
1:B:523:GLN:HA	1:B:523:GLN:NE2	2.24	0.48
1:B:318:GLY:HA2	1:B:492:LEU:HB2	1.94	0.48
1:B:412:LYS:O	1:B:435:VAL:HA	2.12	0.48
1:A:143:VAL:HG21	1:A:236:LEU:HD21	1.96	0.48
1:A:200:ASN:ND2	1:A:200:ASN:N	2.62	0.48
1:B:286:ILE:C	1:B:286:ILE:HD13	2.34	0.48
1:A:400:ILE:HG22	1:A:401:VAL:N	2.27	0.48
1:B:352:LEU:HB3	1:B:412:LYS:HG3	1.95	0.48
1:A:234:ASN:HB2	1:A:236:LEU:HG	1.95	0.48
1:B:39:LYS:HG3	1:B:43:GLY:HA2	1.96	0.48
1:B:147:TRP:CD2	1:B:281:PHE:HE2	2.31	0.48
1:A:19:CYS:HA	1:A:27:LEU:O	2.13	0.48
1:B:195:GLN:HE22	1:B:198:ARG:HD2	1.79	0.48
1:A:50:THR:HG22	3:A:677:HOH:O	2.14	0.48
1:B:131:ASN:HD22	1:B:131:ASN:N	2.12	0.47
1:A:481:LEU:HD13	1:A:484:LEU:HD21	1.95	0.47
1:A:56:PHE:CE1	1:A:461:TYR:HB3	2.49	0.47
1:B:82:VAL:HG21	1:B:154:LEU:CD2	2.44	0.47
1:A:412:LYS:HB3	1:A:412:LYS:NZ	2.25	0.47
1:A:19:CYS:HB3	1:A:28:THR:HA	1.96	0.47
1:A:511:GLU:O	1:A:515:ARG:HG3	2.14	0.47
1:A:533:LEU:CD1	1:B:478:TYR:HD1	2.27	0.47
1:B:18:LYS:O	1:B:28:THR:HA	2.14	0.47
1:A:531:ARG:HD3	1:B:482:THR:HG1	1.73	0.47
1:A:406:LYS:HB3	1:A:407:PRO:HD3	1.97	0.47
1:A:533:LEU:H	1:A:533:LEU:HD22	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:THR:CG2	1:A:297:PRO:HG2	2.45	0.47
1:B:193:ALA:C	1:B:195:GLN:H	2.18	0.47
1:A:445:LEU:C	1:A:445:LEU:HD23	2.34	0.47
1:B:229:ASN:O	1:B:233:GLU:HB3	2.14	0.47
1:B:208:LYS:CB	1:B:210:ASN:HD22	2.28	0.47
1:A:155:TYR:O	1:A:159:GLN:HG3	2.14	0.47
1:A:88:ASN:HD21	1:A:104:ASN:CA	2.28	0.47
1:B:453:THR:HG23	1:B:481:LEU:CD1	2.44	0.47
1:A:67:GLY:HA3	1:A:239:VAL:HG22	1.96	0.47
1:A:154:LEU:HD22	1:A:179:VAL:CG1	2.44	0.47
1:B:26:LEU:HB2	1:B:60:LEU:HD21	1.96	0.47
1:B:300:THR:O	1:B:302:VAL:N	2.45	0.47
1:A:220:LEU:O	1:A:224:ARG:HG3	2.15	0.47
1:B:293:ILE:HG23	1:B:317:ALA:CB	2.45	0.47
1:A:412:LYS:NZ	1:A:414:ALA:HB2	2.29	0.47
1:B:26:LEU:HD12	1:B:27:LEU:H	1.79	0.46
1:B:524:ASN:O	1:B:525:GLU:HB2	2.14	0.46
1:B:266:GLN:HE21	1:B:270:ASN:HD21	1.63	0.46
1:B:270:ASN:O	1:B:271:ASP:HB2	2.15	0.46
1:A:497:PHE:CE2	1:B:530:GLU:HB2	2.50	0.46
1:B:95:GLN:HA	1:B:95:GLN:HE21	1.80	0.46
1:A:350:ASN:HD22	1:A:412:LYS:HE2	1.80	0.46
1:B:353:GLY:HA3	1:B:409:GLY:O	2.16	0.46
1:A:428:ASN:ND2	1:B:436:CYS:HB3	2.30	0.46
1:B:153:ASP:OD2	1:B:155:TYR:N	2.42	0.46
1:A:378:ASP:HA	1:A:381:ILE:CG2	2.46	0.46
1:B:258:ASN:ND2	1:B:258:ASN:N	2.63	0.46
1:A:153:ASP:C	1:A:153:ASP:OD2	2.54	0.46
1:B:380:ILE:O	1:B:380:ILE:HG22	2.16	0.46
1:A:256:GLY:HA2	1:A:263:ASN:ND2	2.31	0.46
1:B:134:LEU:HB3	1:B:135:PRO:CD	2.47	0.45
1:A:286:ILE:HA	1:A:314:THR:HG21	1.96	0.45
1:B:76:ASN:HB3	1:B:439:SER:OG	2.16	0.45
1:B:131:ASN:N	1:B:131:ASN:ND2	2.64	0.45
1:B:122:GLU:N	1:B:122:GLU:CD	2.70	0.45
1:B:465:ASP:OD2	1:B:467:VAL:HB	2.16	0.45
1:A:14:VAL:HG22	1:B:47:VAL:CG1	2.47	0.45
1:B:97:LYS:HB2	1:B:97:LYS:HZ2	1.80	0.45
1:A:55:VAL:HG23	1:A:464:VAL:HG21	1.99	0.45
1:A:335:PHE:CZ	1:B:327:LYS:HE2	2.51	0.45
1:A:258:ASN:ND2	1:A:258:ASN:N	2.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:TYR:CZ	1:B:188:PRO:HB3	2.52	0.45
1:A:223:ILE:O	1:A:226:ASP:HB2	2.17	0.45
1:B:389:ASN:OD1	1:B:392:LEU:HD13	2.17	0.45
1:B:192:ALA:C	1:B:194:ASN:N	2.68	0.45
1:B:391:LYS:HB2	1:B:391:LYS:NZ	2.32	0.45
1:A:44:ARG:HH21	1:B:13:LYS:CG	2.27	0.45
1:B:200:ASN:HD22	1:B:200:ASN:H	1.65	0.45
1:A:337:VAL:HG21	1:A:380:ILE:HG22	1.98	0.45
1:B:130:PHE:CE2	1:B:136:MET:HE1	2.51	0.45
1:A:352:LEU:CD2	1:A:352:LEU:N	2.79	0.45
1:A:122:GLU:OE2	1:A:122:GLU:N	2.50	0.45
1:A:325:GLN:HG3	3:A:652:HOH:O	2.16	0.44
1:A:183:PRO:HB2	1:A:203:ILE:HG23	2.00	0.44
1:B:150:ASN:ND2	1:B:160:ARG:NH1	2.64	0.44
1:A:378:ASP:CA	1:A:381:ILE:HG22	2.47	0.44
1:A:286:ILE:HG23	1:A:314:THR:HB	1.98	0.44
1:A:115:LEU:HD22	1:A:511:GLU:HB3	2.00	0.44
1:B:83:ALA:HB2	1:B:176:MET:HE2	1.99	0.44
1:A:351:HIS:ND1	1:A:351:HIS:O	2.51	0.44
1:B:302:VAL:HG23	1:B:305:LEU:H	1.82	0.44
1:B:285:SER:HB3	1:B:290:VAL:HG23	1.99	0.44
1:B:26:LEU:HD12	1:B:27:LEU:N	2.33	0.44
1:B:352:LEU:N	1:B:352:LEU:CD2	2.81	0.44
1:B:272:HIS:ND1	1:B:274:GLU:N	2.59	0.44
1:A:301:PHE:HA	1:A:305:LEU:HD23	1.98	0.44
1:B:293:ILE:CA	1:B:317:ALA:HB2	2.35	0.44
1:A:465:ASP:OD1	1:A:467:VAL:HG22	2.18	0.44
1:A:500:VAL:HG21	1:B:527:ARG:NH2	2.32	0.44
1:B:438:ASP:HB3	2:B:660:NAD:O7N	2.18	0.44
1:B:246:ASN:HD22	2:B:660:NAD:H8A	1.81	0.44
1:A:57:LYS:HB3	1:A:460:SER:HB3	1.98	0.44
1:A:169:GLN:HB3	1:A:169:GLN:HE21	1.61	0.44
1:B:293:ILE:HA	1:B:317:ALA:HB3	1.93	0.44
1:B:174:ALA:HB3	3:B:669:HOH:O	2.18	0.44
1:B:350:ASN:O	1:B:351:HIS:HB3	2.18	0.44
1:B:259:ASP:OD2	1:B:259:ASP:N	2.49	0.44
1:B:348:SER:OG	1:B:416:ASP:HA	2.18	0.44
1:A:52:GLN:OE1	1:A:53:ASP:O	2.36	0.44
1:B:254:SER:H	1:B:258:ASN:ND2	2.03	0.43
1:B:468:LYS:HG2	1:B:471:ALA:CB	2.48	0.43
1:A:530:GLU:HG3	1:B:497:PHE:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:VAL:HB	1:A:303:PRO:HD2	2.00	0.43
1:A:109:MET:HE2	1:A:507:ARG:HE	1.83	0.43
1:A:104:ASN:HD22	1:A:104:ASN:C	2.21	0.43
1:A:184:SER:OG	1:A:185:ILE:N	2.49	0.43
1:B:94:PHE:CD2	1:B:94:PHE:N	2.86	0.43
1:B:254:SER:O	1:B:257:VAL:HG13	2.18	0.43
1:A:532:LEU:HG	1:B:461:TYR:CZ	2.54	0.43
1:A:520:LEU:HA	1:A:520:LEU:HD23	1.77	0.43
1:A:525:GLU:OE2	1:A:525:GLU:HA	2.18	0.43
1:A:440:LEU:HA	1:A:440:LEU:HD12	1.83	0.43
1:A:447:ILE:O	1:A:451:VAL:HG23	2.19	0.43
1:A:38:THR:O	1:A:45:PHE:HA	2.19	0.43
1:A:51:VAL:HG12	1:A:52:GLN:N	2.34	0.43
1:A:125:ASP:OD1	1:A:527:ARG:NH2	2.51	0.43
1:A:47:VAL:HG11	1:B:520:LEU:HD11	2.01	0.43
1:A:190:PHE:HB3	1:A:248:GLU:HB3	2.01	0.43
1:A:134:LEU:HD21	1:A:518:ILE:CG2	2.48	0.43
1:B:260:THR:OG1	1:B:263:ASN:ND2	2.52	0.43
1:B:115:LEU:O	1:B:127:TYR:HA	2.19	0.43
1:A:354:ASN:HD21	1:A:356:ASP:HB2	1.83	0.43
1:A:234:ASN:CB	1:A:236:LEU:HG	2.49	0.43
1:B:350:ASN:HD21	1:B:402:ILE:CG1	2.32	0.42
1:B:526:LEU:O	1:B:527:ARG:C	2.57	0.42
1:B:276:ALA:O	1:B:279:THR:HB	2.19	0.42
1:B:381:ILE:HG23	1:B:388:TYR:CG	2.54	0.42
1:A:533:LEU:HD22	1:A:533:LEU:N	2.34	0.42
1:B:184:SER:OG	1:B:185:ILE:N	2.52	0.42
1:A:14:VAL:HG22	1:B:47:VAL:HG11	2.00	0.42
1:A:483:PHE:CE1	1:B:531:ARG:HD3	2.55	0.42
1:A:349:TYR:HD1	1:A:401:VAL:HB	1.84	0.42
1:B:350:ASN:ND2	1:B:402:ILE:HG13	2.35	0.42
1:A:476:ASN:O	1:A:477:PHE:C	2.58	0.42
1:B:501:ASN:HA	1:B:506:GLN:OE1	2.19	0.42
1:B:163:VAL:HG23	1:B:164:LEU:HD13	2.01	0.42
1:B:473:LYS:CD	1:B:474:PHE:H	2.30	0.42
1:A:247:THR:HG23	1:A:298:GLN:HE21	1.84	0.42
1:B:71:ILE:HG13	1:B:147:TRP:CE3	2.54	0.42
1:A:242:LEU:HD12	1:A:293:ILE:HD12	2.01	0.42
1:A:250:TYR:CD2	1:A:299:ASN:OD1	2.73	0.42
1:A:10:THR:HG21	1:A:129:PRO:HG2	2.00	0.42
1:B:258:ASN:H	1:B:258:ASN:ND2	2.13	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:GLU:OE1	1:B:277:PRO:HD2	2.20	0.42
1:A:516:LEU:C	1:A:516:LEU:CD2	2.88	0.42
1:B:195:GLN:HA	1:B:195:GLN:NE2	2.35	0.42
1:B:384:ASN:ND2	1:B:387:LEU:HB2	2.35	0.42
1:B:525:GLU:OE2	1:B:527:ARG:NH1	2.53	0.42
1:B:265:LEU:O	1:B:269:LYS:HG3	2.19	0.42
1:B:322:LYS:NZ	1:B:503:LEU:HD13	2.35	0.42
1:B:493:THR:CG2	1:B:499:PRO:HB3	2.50	0.42
1:B:184:SER:O	1:B:202:CYS:HA	2.20	0.42
1:B:185:ILE:CG1	1:B:203:ILE:HD11	2.48	0.42
1:A:294:ASN:ND2	1:A:294:ASN:C	2.72	0.42
1:B:56:PHE:HA	1:B:461:TYR:HA	2.02	0.42
1:A:19:CYS:CB	1:A:28:THR:OG1	2.67	0.42
1:A:264:LEU:HD21	1:A:305:LEU:HD13	2.00	0.42
1:A:328:LEU:HD13	1:B:335:PHE:CE1	2.55	0.41
1:A:347:ALA:O	1:A:416:ASP:HA	2.20	0.41
1:B:95:GLN:CA	1:B:95:GLN:NE2	2.83	0.41
1:B:208:LYS:HB2	1:B:210:ASN:HD22	1.85	0.41
1:A:80:THR:HG23	1:A:164:LEU:HD11	2.01	0.41
1:A:129:PRO:HD2	1:A:132:SER:OG	2.20	0.41
1:A:59:ASP:O	1:A:458:ARG:HD2	2.20	0.41
1:B:498:HIS:CD2	1:B:498:HIS:N	2.87	0.41
1:B:468:LYS:HZ3	1:B:471:ALA:CB	2.34	0.41
1:A:412:LYS:HG3	1:A:438:ASP:OD1	2.20	0.41
1:B:71:ILE:HG21	1:B:243:TRP:CZ3	2.56	0.41
1:A:73:LEU:HD23	1:A:181:PRO:HB3	2.01	0.41
1:A:325:GLN:HB3	1:A:325:GLN:HE21	1.68	0.41
1:A:150:ASN:ND2	1:A:151:ASN:N	2.68	0.41
1:B:231:LYS:HG2	1:B:239:VAL:HG21	2.03	0.41
1:A:56:PHE:CD1	1:A:461:TYR:HB3	2.55	0.41
1:B:268:ILE:HG12	1:B:275:ILE:HD12	2.03	0.41
1:A:533:LEU:HD23	1:B:461:TYR:OH	2.20	0.41
1:B:148:ASP:HA	2:B:660:NAD:N3A	2.35	0.41
1:A:67:GLY:CA	1:A:236:LEU:HD13	2.50	0.41
1:A:10:THR:HG23	1:A:132:SER:CB	2.50	0.41
1:A:225:ARG:HD2	1:A:225:ARG:HA	1.85	0.41
1:A:240:ILE:HG23	1:A:291:PRO:HB2	2.02	0.41
1:A:104:ASN:C	1:A:104:ASN:ND2	2.74	0.41
1:B:70:LEU:HD12	1:B:242:LEU:HB3	2.02	0.41
1:A:406:LYS:N	1:A:407:PRO:CD	2.83	0.41
1:A:320:ASP:HB3	1:A:489:LYS:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LYS:HA	1:A:175:LYS:HD3	1.83	0.41
1:A:328:LEU:HD11	1:B:332:LEU:HG	2.02	0.41
1:B:254:SER:O	1:B:258:ASN:ND2	2.54	0.41
1:B:486:TYR:CE1	1:B:507:ARG:HB2	2.56	0.41
1:B:195:GLN:NE2	1:B:198:ARG:HD2	2.36	0.41
1:B:424:LEU:HA	3:B:702:HOH:O	2.19	0.41
1:A:215:GLY:HA2	3:A:660:HOH:O	2.21	0.41
1:B:323:SER:O	1:B:327:LYS:HG3	2.21	0.41
1:B:349:TYR:N	1:B:349:TYR:CD1	2.88	0.41
1:A:322:LYS:HD2	1:A:486:TYR:O	2.21	0.41
1:A:354:ASN:HD21	1:A:356:ASP:CB	2.34	0.41
1:A:532:LEU:HD23	1:B:54:TYR:CE2	2.55	0.41
1:A:21:TYR:CZ	1:A:26:LEU:HD13	2.56	0.41
1:A:381:ILE:HG13	1:A:388:TYR:CD2	2.55	0.41
1:A:247:THR:HG23	1:A:298:GLN:NE2	2.36	0.41
1:A:507:ARG:O	1:A:510:LEU:HB2	2.21	0.41
1:B:119:ILE:HA	1:B:124:ASN:O	2.21	0.41
1:A:174:ALA:HB3	3:A:709:HOH:O	2.20	0.41
1:A:228:GLN:HG2	1:A:288:GLU:OE1	2.21	0.40
1:A:195:GLN:NE2	1:A:355:ASN:HD21	2.15	0.40
1:A:452:MET:HG3	1:A:487:TRP:HH2	1.86	0.40
1:B:153:ASP:OD2	1:B:155:TYR:HB3	2.22	0.40
1:A:10:THR:HG23	1:A:132:SER:HB2	2.02	0.40
1:A:200:ASN:HD22	1:A:200:ASN:C	2.23	0.40
1:A:26:LEU:CB	1:A:60:LEU:HD11	2.52	0.40
1:B:245:ALA:HA	2:B:660:NAD:H52A	2.02	0.40
1:B:483:PHE:HA	1:B:483:PHE:HD2	1.81	0.40
1:A:71:ILE:HG21	1:A:243:TRP:CZ3	2.56	0.40
1:B:322:LYS:HD2	1:B:489:LYS:HA	2.03	0.40
1:B:507:ARG:NH2	1:B:507:ARG:HG3	2.30	0.40
1:A:418:TYR:N	1:A:418:TYR:CD1	2.90	0.40
1:B:477:PHE:O	1:B:478:TYR:C	2.59	0.40
1:B:484:LEU:HA	1:B:487:TRP:CZ3	2.57	0.40
1:A:482:THR:OG1	1:B:531:ARG:HG2	2.21	0.40
1:A:526:LEU:O	1:A:531:ARG:HG3	2.21	0.40
1:B:208:LYS:HB3	1:B:210:ASN:ND2	2.37	0.40
1:A:109:MET:HE2	1:A:507:ARG:NE	2.36	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	502/533 (94%)	431 (86%)	57 (11%)	14 (3%)	6	5
1	B	494/533 (93%)	429 (87%)	52 (10%)	13 (3%)	7	6
All	All	996/1066 (93%)	860 (86%)	109 (11%)	27 (3%)	6	6

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	SER
1	B	199	ALA
1	B	319	ASP
1	B	320	ASP
1	B	468	LYS
1	A	194	ASN
1	A	196	ASP
1	A	206	ASP
1	A	209	GLY
1	A	324	GLY
1	A	351	HIS
1	B	133	LEU
1	B	193	ALA
1	B	204	ASN
1	B	301	PHE
1	B	191	ILE
1	A	319	ASP
1	A	407	PRO
1	A	470	ASP
1	B	22	LYS
1	B	495	PRO
1	A	530	GLU
1	B	346	ILE
1	A	346	ILE
1	A	401	VAL

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Mol	Chain	Res	Type
1	A	480	VAL
1	B	479	PRO

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	447/471 (95%)	409 (92%)	38 (8%)	13	20
1	B	439/471 (93%)	399 (91%)	40 (9%)	12	17
All	All	886/942 (94%)	808 (91%)	78 (9%)	12	18

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU
1	A	48	THR
1	A	58	LEU
1	A	66	LEU
1	A	86	LEU
1	A	104	ASN
1	A	110	THR
1	A	111	GLN
1	A	122	GLU
1	A	150	ASN
1	A	189	ASP
1	A	200	ASN
1	A	201	ASN
1	A	207	GLU
1	A	211	VAL
1	A	249	ARG
1	A	258	ASN
1	A	260	THR
1	A	261	MET
1	A	262	GLU
1	A	298	GLN

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Mol	Chain	Res	Type
1	A	300	THR
1	A	320	ASP
1	A	332	LEU
1	A	336	LEU
1	A	352	LEU
1	A	355	ASN
1	A	378	ASP
1	A	390	ASP
1	A	397	ASP
1	A	401	VAL
1	A	412	LYS
1	A	418	TYR
1	A	428	ASN
1	A	438	ASP
1	A	473	LYS
1	A	516	LEU
1	A	532	LEU
1	B	70	LEU
1	B	73	LEU
1	B	81	LEU
1	B	92	VAL
1	B	111	GLN
1	B	116	LYS
1	B	131	ASN
1	B	151	ASN
1	B	163	VAL
1	B	164	LEU
1	B	168	LEU
1	B	196	ASP
1	B	197	GLU
1	B	200	ASN
1	B	207	GLU
1	B	214	ARG
1	B	249	ARG
1	B	250	TYR
1	B	254	SER
1	B	258	ASN
1	B	262	GLU
1	B	286	ILE
1	B	299	ASN
1	B	308	LEU
1	B	335	PHE

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Mol	Chain	Res	Type
1	B	349	TYR
1	B	352	LEU
1	B	387	LEU
1	B	394	LYS
1	B	397	ASP
1	B	408	VAL
1	B	469	GLU
1	B	475	GLU
1	B	476	ASN
1	B	479	PRO
1	B	483	PHE
1	B	488	LEU
1	B	507	ARG
1	B	517	LEU
1	B	523	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (55) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	77	ASN
1	A	88	ASN
1	A	91	ASN
1	A	104	ASN
1	A	111	GLN
1	A	150	ASN
1	A	151	ASN
1	A	162	GLN
1	A	169	GLN
1	A	170	GLN
1	A	200	ASN
1	A	201	ASN
1	A	210	ASN
1	A	228	GLN
1	A	246	ASN
1	A	258	ASN
1	A	263	ASN
1	A	270	ASN
1	A	272	HIS
1	A	294	ASN
1	A	298	GLN
1	A	307	GLN

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Mol	Chain	Res	Type
1	A	325	GLN
1	A	355	ASN
1	A	384	ASN
1	A	428	ASN
1	A	476	ASN
1	A	504	ASN
1	A	512	ASN
1	A	523	GLN
1	B	91	ASN
1	B	95	GLN
1	B	131	ASN
1	B	150	ASN
1	B	151	ASN
1	B	159	GLN
1	B	170	GLN
1	B	195	GLN
1	B	200	ASN
1	B	201	ASN
1	B	210	ASN
1	B	219	HIS
1	B	221	GLN
1	B	228	GLN
1	B	246	ASN
1	B	258	ASN
1	B	263	ASN
1	B	270	ASN
1	B	299	ASN
1	B	307	GLN
1	B	350	ASN
1	B	476	ASN
1	B	506	GLN
1	B	523	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

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	$\# Z > 2$
2	NAD	A	650	-	38,48,48	1.73	4 (10%)	47,73,73	1.34	8 (17%)
2	NAD	B	660	-	38,48,48	1.87	7 (18%)	47,73,73	1.44	6 (12%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAD	A	650	-	-	0/22/62/62	0/5/5/5
2	NAD	B	660	-	-	0/22/62/62	0/5/5/5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	660	NAD	O4B-C1B	-4.27	1.35	1.41
2	B	660	NAD	C2N-C3N	2.41	1.42	1.39
2	B	660	NAD	C3N-C7N	2.51	1.54	1.50
2	B	660	NAD	C5N-C4N	3.37	1.45	1.38
2	A	650	NAD	C5N-C4N	3.65	1.46	1.38
2	A	650	NAD	C2A-N1A	3.79	1.41	1.33
2	A	650	NAD	C6N-N1N	3.85	1.45	1.35
2	B	660	NAD	C2A-N1A	3.92	1.41	1.33
2	B	660	NAD	C6N-N1N	3.94	1.46	1.35
2	A	650	NAD	C4N-C3N	6.15	1.49	1.39
2	B	660	NAD	C4N-C3N	6.37	1.50	1.39

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	650	NAD	C1B-N9A-C4A	-3.34	121.90	126.94
2	A	650	NAD	N3A-C2A-N1A	-2.65	126.86	128.89
2	B	660	NAD	O7N-C7N-N7N	-2.53	119.03	122.59
2	B	660	NAD	C1B-N9A-C4A	-2.41	123.31	126.94
2	A	650	NAD	O7N-C7N-N7N	-2.33	119.32	122.59
2	B	660	NAD	O2A-PA-O3	2.10	114.64	105.09
2	A	650	NAD	O2A-PA-O3	2.25	115.31	105.09
2	A	650	NAD	O4D-C4D-C3D	2.36	109.89	105.15
2	B	660	NAD	C4A-C5A-N7A	2.49	111.77	109.48
2	A	650	NAD	C4A-C5A-N7A	2.51	111.79	109.48
2	A	650	NAD	C3N-C7N-N7N	3.18	121.30	117.82
2	B	660	NAD	C3N-C7N-N7N	3.44	121.58	117.82
2	A	650	NAD	O4D-C1D-N1N	4.30	112.86	108.13
2	B	660	NAD	O4D-C1D-N1N	5.82	114.53	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	650	NAD	2	0
2	B	660	NAD	5	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	506/533 (94%)	0.61	56 (11%) 7 7	26, 53, 91, 91	0
1	B	498/533 (93%)	0.85	72 (14%) 3 3	27, 56, 91, 91	0
All	All	1004/1066 (94%)	0.73	128 (12%) 5 5	26, 54, 91, 91	0

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	467	VAL	10.3
1	B	404	TYR	9.0
1	B	392	LEU	8.6
1	B	402	ILE	7.6
1	B	399	CYS	7.5
1	A	392	LEU	7.4
1	A	405	MET	7.3
1	A	381	ILE	7.2
1	B	469	GLU	7.0
1	A	404	TYR	6.8
1	B	400	ILE	6.8
1	A	193	ALA	6.8
1	B	468	LYS	6.8
1	B	394	LYS	6.6
1	B	393	GLY	6.4
1	A	194	ASN	6.4
1	B	348	SER	6.4
1	A	396	VAL	6.2
1	B	466	PRO	6.1
1	A	400	ILE	5.8
1	A	377	ILE	5.4
1	A	9	ILE	5.3
1	B	380	ILE	5.2
1	B	390	ASP	5.2

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Mol	Chain	Res	Type	RSRZ
1	A	391	LYS	5.2
1	A	378	ASP	5.1
1	B	382	ALA	5.0
1	A	402	ILE	5.0
1	B	391	LYS	4.9
1	B	193	ALA	4.9
1	B	396	VAL	4.9
1	B	317	ALA	4.8
1	A	393	GLY	4.8
1	B	194	ASN	4.8
1	B	389	ASN	4.7
1	A	390	ASP	4.7
1	B	471	ALA	4.7
1	B	352	LEU	4.6
1	B	497	PHE	4.6
1	A	406	LYS	4.6
1	A	398	HIS	4.4
1	B	500	VAL	4.4
1	B	385	ASP	4.3
1	A	382	ALA	4.3
1	A	385	ASP	4.3
1	B	470	ASP	4.2
1	B	351	HIS	4.2
1	B	403	LYS	4.1
1	A	380	ILE	4.1
1	B	349	TYR	4.1
1	A	379	ASP	4.0
1	A	192	ALA	4.0
1	B	398	HIS	4.0
1	B	381	ILE	4.0
1	B	397	ASP	3.9
1	A	376	VAL	3.9
1	A	40	THR	3.9
1	B	408	VAL	3.9
1	B	499	PRO	3.8
1	A	394	LYS	3.8
1	B	386	ILE	3.7
1	B	410	ASP	3.6
1	A	41	ALA	3.5
1	B	347	ALA	3.5
1	A	397	ASP	3.5
1	B	405	MET	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	197	GLU	3.5
1	B	395	LYS	3.5
1	A	208	LYS	3.5
1	A	386	ILE	3.4
1	B	40	THR	3.3
1	A	407	PRO	3.3
1	B	21	TYR	3.2
1	A	408	VAL	3.2
1	B	409	GLY	3.2
1	A	399	CYS	3.2
1	B	465	ASP	3.2
1	A	389	ASN	3.2
1	B	383	SER	3.2
1	B	23	ASP	3.1
1	B	407	PRO	3.1
1	A	395	LYS	3.0
1	A	388	TYR	2.9
1	B	192	ALA	2.9
1	A	195	GLN	2.9
1	B	498	HIS	2.9
1	A	355	ASN	2.9
1	B	384	ASN	2.8
1	A	197	GLU	2.7
1	B	495	PRO	2.7
1	B	350	ASN	2.7
1	B	388	TYR	2.6
1	B	242	LEU	2.6
1	B	208	LYS	2.6
1	A	259	ASP	2.6
1	A	15	VAL	2.5
1	A	352	LEU	2.5
1	A	401	VAL	2.5
1	B	406	LYS	2.5
1	B	353	GLY	2.5
1	B	41	ALA	2.5
1	B	401	VAL	2.4
1	A	533	LEU	2.4
1	B	533	LEU	2.4
1	B	472	GLY	2.4
1	B	45	PHE	2.4
1	A	460	SER	2.4
1	B	474	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	412	LYS	2.3
1	A	42	SER	2.3
1	A	351	HIS	2.3
1	B	346	ILE	2.3
1	A	44	ARG	2.3
1	A	470	ASP	2.3
1	A	123	GLY	2.2
1	A	497	PHE	2.2
1	B	293	ILE	2.2
1	B	488	LEU	2.2
1	A	500	VAL	2.1
1	B	387	LEU	2.1
1	B	43	GLY	2.1
1	A	383	SER	2.1
1	B	241	VAL	2.1
1	B	473	LYS	2.1
1	A	207	GLU	2.1
1	B	411	SER	2.1
1	A	387	LEU	2.0
1	A	440	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAD	B	660	44/44	0.95	0.20	0.11	43,56,79,81	0
2	NAD	A	650	44/44	0.95	0.18	-0.02	36,57,84,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.