



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:17 PM UTC

PDB ID : 1P1I / pdb_00001p1i
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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

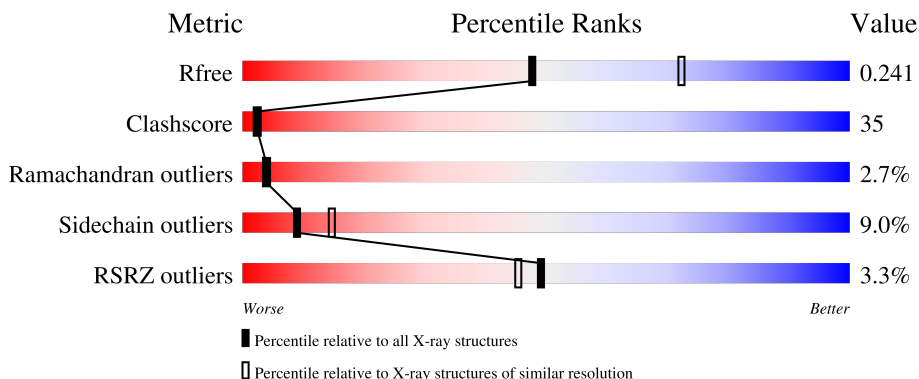
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

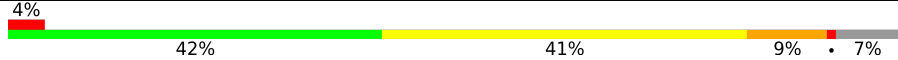
The reported resolution of this entry is 2.40 Å.

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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 $\leq 5\%$. 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	
1	B	533	

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.

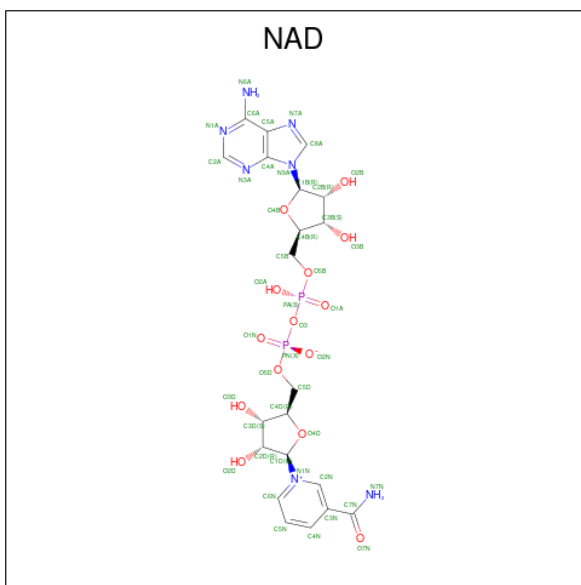
- Molecule 1 is a protein called Inositol-3-phosphate synthase.

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 (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is water.

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

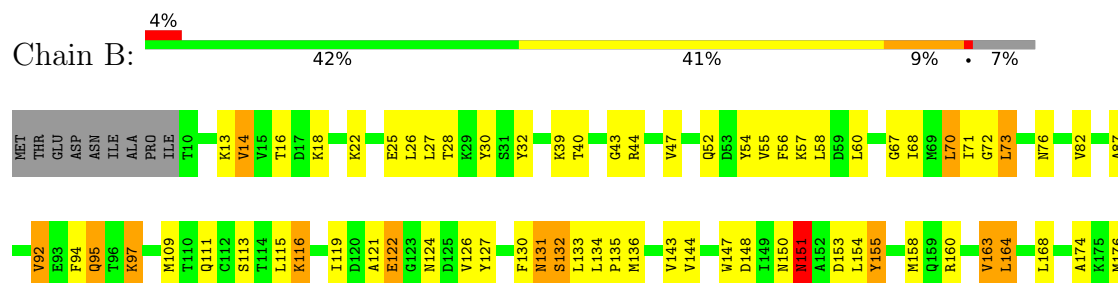
3 Residue-property plots

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 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	152.38Å 97.34Å 122.89Å 90.00° 126.53° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 10.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.40) 98.2 (10.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.209 , 0.288 0.246 , 0.241	Depositor DCC
R_{free} test set	2857 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	46.4	Xtriage
Anisotropy	0.473	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 80.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.004 for h,-k,-h-l	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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.47	0/4066	0.97	10/5514 (0.2%)
1	B	0.48	0/4003	0.98	18/5427 (0.3%)
All	All	0.47	0/8069	0.98	28/10941 (0.3%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	GLY	N-CA-C	9.70	124.91	110.75
1	A	377	ILE	N-CA-C	-8.79	104.91	113.53
1	B	155	TYR	N-CA-C	-8.51	101.59	111.03
1	B	494	ARG	N-CA-C	-7.83	98.26	109.62
1	B	208	LYS	N-CA-C	-7.64	104.42	113.21
1	A	478	TYR	N-CA-C	-7.43	101.06	109.60
1	B	391	LYS	N-CA-C	-6.93	104.81	113.55
1	B	319	ASP	N-CA-C	6.78	125.25	110.80
1	B	194	ASN	N-CA-C	-6.68	105.44	112.93
1	B	151	ASN	N-CA-C	6.51	120.14	112.72
1	B	406	LYS	CA-C-N	6.45	126.20	119.05
1	B	406	LYS	C-N-CA	6.45	126.20	119.05
1	B	144	VAL	N-CA-C	6.25	117.59	108.53
1	B	132	SER	N-CA-C	6.24	118.89	111.71
1	A	464	VAL	N-CA-C	5.99	116.50	108.11
1	B	381	ILE	N-CA-C	-5.59	105.47	113.07
1	A	485	SER	N-CA-C	5.58	118.12	111.71
1	A	193	ALA	N-CA-C	5.55	117.78	111.11
1	A	143	VAL	N-CA-C	-5.50	98.77	106.53
1	B	143	VAL	N-CA-C	-5.46	98.60	107.28
1	B	210	ASN	N-CA-C	5.42	117.29	110.24
1	A	151	ASN	N-CA-C	5.41	119.59	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	144	VAL	N-CA-C	5.28	116.90	108.87
1	B	72	GLY	N-CA-C	-5.18	107.43	115.67
1	B	466	PRO	N-CA-C	-5.17	107.89	114.35
1	A	323	SER	N-CA-C	5.14	117.48	109.52
1	B	410	ASP	N-CA-C	-5.10	106.90	113.23
1	B	387	LEU	N-CA-C	-5.03	107.70	113.88

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3989	0	3994	283	0
1	B	3926	0	3939	312	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	554	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 35.

All (554) 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:109:MET:HE2	1:B:507:ARG:NH2	1.85	0.91
1:B:476:ASN:H	1:B:476:ASN:HD22	1.15	0.90
1:A:104:ASN:ND2	1:A:106:PHE:H	1.69	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:37:VAL:HG11	1:B:126:VAL:HG11	1.61	0.81
1:A:104:ASN:HD22	1:A:106:PHE:H	1.27	0.81
1:B:58:LEU:HD23	1:B:459:VAL:HG22	1.64	0.80
1:A:158:MET:HE2	1:A:176:MET:HE2	1.61	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	0.76
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:B:238:LYS:NZ	1:B:457:THR:HG21	2.01	0.75
1:B:115:LEU:HD22	1:B:511:GLU:HG3	1.67	0.75
1:B:135:PRO:HG2	1:B:458:ARG:HH22	1.51	0.75
1:A:423:MET:HE3	1:A:424:LEU:HG	1.67	0.74
1:B:286:ILE:HD11	1:B:308:LEU:HD22	1.68	0.74
1:A:352:LEU:HD12	1:A:354:ASN:OD1	1.88	0.73
1:B:286:ILE:CD1	1:B:308:LEU:HD13	2.17	0.73
1:B:404:TYR:O	1:B:405:MET:HE2	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:391:LYS:HB2	1:B:391:LYS:HZ2	1.54	0.72
1:B:261:MET:HE1	1:B:311:HIS:HB2	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HD23	1:A:352:LEU:H	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:150:ASN:HD21	1:B:160:ARG:HH12	1.38	0.71
1:B:476:ASN:HD22	1:B:476:ASN:N	1.84	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: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:A:150:ASN:ND2	1:A:152:ALA:H	1.89	0.69
1:B:254:SER:N	1:B:258:ASN:HD21	1.90	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:A:353:GLY:CA	1:A:406:LYS:HA	2.23	0.68
1:B:294:ASN:H	1:B:317:ALA:HB3	1.58	0.68
1:A:158:MET:HE2	1:A:176:MET:HG3	1.74	0.68
1:B:131:ASN:HD22	1:B:132:SER:N	1.93	0.67
1:B:454:GLU:O	1:B:457:THR:HB	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:B:323:SER:HB3	1:B:489:LYS:HZ1	1.59	0.67
1:A:405:MET:HE2	1:A:405:MET:HA	1.76	0.67
1:A:272:HIS:CD2	1:A:274:GLU:HB2	2.29	0.67
1:A:400:ILE:HG22	1:A:401:VAL:H	1.58	0.67
1:A:473:LYS:H	1:A:473:LYS:HD2	1.60	0.67
1:A:352:LEU:HD23	1:A:352:LEU:N	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
1:B:384:ASN:ND2	1:B:387:LEU:HD22	2.10	0.66
1:A:246:ASN:ND2	2:A:650:NAD:H8A	2.05	0.66
1:A:328:LEU:HD21	1:B:430:ILE:HG21	1.78	0.66
1:B:507:ARG:HH21	1:B:507:ARG:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:109:MET:HE2	1:B:507:ARG:HH21	1.61	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:A:481:LEU:HB3	1:A:484:LEU:CD2	2.28	0.64
1:B:203:ILE:HG13	1:B:204:ASN:N	2.13	0.64
1:A:478:TYR:CD2	1:A:479:PRO:HD2	2.32	0.64
1:B:478:TYR:CD1	1:B:479:PRO:HD2	2.33	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:A:190:PHE:O	1:A:248:GLU:HA	1.98	0.63
1:B:121:ALA:HB3	1:B:122:GLU:OE2	1.97	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:277:PRO:HG2	3:B:665:HOH:O	1.99	0.62
1:B:387:LEU:HA	1:B:392:LEU:HD22	1.80	0.62
1:B:249:ARG:HG2	1:B:249:ARG:HH21	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:A:516:LEU:C	1:A:516:LEU:HD23	2.24	0.61
1:B:30:TYR:CE1	1:B:32:TYR:HB2	2.35	0.61
1:B:40:THR:HG23	1:B:44:ARG:O	1.99	0.61
1:B:352:LEU:N	1:B:352:LEU:HD23	2.15	0.61
1:B:261:MET:H	1:B:307:GLN:NE2	1.99	0.61
1:A:406:LYS:HD3	1:A:406:LYS:C	2.25	0.61
1:B:200:ASN:C	1:B:200:ASN:ND2	2.58	0.61
1:B:515:ARG:NH1	1:B:521:PRO:O	2.33	0.61
1:A:14:VAL:HA	1:B:47:VAL:HG13	1.81	0.61
1:B:57:LYS:HB2	1:B:474:PHE:CE2	2.35	0.61
1:A:125:ASP:CG	1:A:527:ARG:HH22	2.07	0.61
1:A:400:ILE:O	1:A:401:VAL:HG23	2.00	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:B:56:PHE:CD2	1:B:461:TYR:HB3	2.34	0.61
1:B:255:PRO:HA	1:B:259:ASP:CG	2.25	0.61
1:B:302:VAL:HB	1:B:303:PRO:HD2	1.83	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:381:ILE:HG13	1:A:388:TYR:CG	2.36	0.60
1:A:486:TYR:HA	1:A:506:GLN:OE1	2.01	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:217:TRP:O	1:A:221:GLN:HG2	2.01	0.60
1:A:272:HIS:CD2	1:A:274:GLU:H	2.20	0.60
1:B:220:LEU:C	1:B:220:LEU:HD23	2.27	0.59
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:A:19:CYS:HB3	1:A:28:THR:OG1	2.03	0.59
1:B:469:GLU:HG2	1:B:470:ASP:N	2.17	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:350:ASN:O	1:A:351:HIS:HB3	2.02	0.59
1:A:509:ALA:CB	1:B:531:ARG:HH12	2.13	0.59
1:B:332:LEU:HD23	1:B:332:LEU:C	2.28	0.59
1:B:349:TYR:CB	1:B:401:VAL:HB	2.31	0.59
1:A:158:MET:HE2	1:A:176:MET:CE	2.31	0.59
1:A:158:MET:CE	1:A:176:MET:HG3	2.32	0.59
1:B:261:MET:HE1	1:B:311:HIS:CB	2.33	0.59
1:A:38:THR:HG22	1:A:39:LYS:N	2.18	0.59
1:B:200:ASN:C	1:B:200:ASN:HD22	2.09	0.58
1:A:294:ASN:ND2	1:A:296:SER:H	2.01	0.58
1:B:238:LYS:HZ1	1:B:457:THR:HG21	1.68	0.58
1:B:352:LEU:HD21	1:B:402:ILE:HD11	1.86	0.58
1:B:516:LEU:HD12	1:B:516:LEU:C	2.29	0.58
1:A:14:VAL:HG21	1:A:518:ILE:HD12	1.85	0.58
1:B:299:ASN:HD22	1:B:299:ASN:H	1.52	0.58
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:HG12	1:A:92:VAL:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:PRO:HB3	1:A:320:ASP:OD2	2.04	0.57
1:B:134:LEU:HD21	1:B:518:ILE:CG2	2.34	0.57
1:A:44:ARG:NH2	1:B:13:LYS:HG2	2.20	0.57
1:B:58:LEU:HD12	1:B:60:LEU:HD23	1.86	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:332:LEU:HD13	1:B:432:ILE:CD1	2.16	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:208:LYS:C	1:B:210:ASN:H	2.13	0.56
1:B:26:LEU:O	1:B:57:LYS:HA	2.04	0.56
1:B:489:LYS:O	1:B:491:PRO:HD3	2.05	0.56
1:B:97:LYS:HB2	1:B:97:LYS:NZ	2.20	0.56
1:B:134:LEU:HD12	1:B:455:PHE:HZ	1.70	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: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:30:TYR:HE1	1:B:32:TYR:HB2	1.71	0.56
1:B:310:GLU:HG2	1:B:479:PRO:HG2	1.87	0.56
1:A:494:ARG:HB2	1:A:495:PRO:HD2	1.88	0.55
1:A:528:PHE:C	1:A:530:GLU:H	2.14	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:476:ASN:ND2	1:B:476:ASN:N	2.47	0.55
1:A:200:ASN:HD22	1:A:200:ASN:N	2.02	0.55
1:A:39:LYS:HA	1:A:44:ARG:O	2.05	0.55
1:B:207:GLU:C	1:B:209:GLY:H	2.13	0.55
1:A:256:GLY:C	1:A:263:ASN:HD22	2.15	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:A:150:ASN:HD22	1:A:150:ASN:C	2.15	0.55
1:B:468:LYS:NZ	1:B:471:ALA:CB	2.69	0.55
1:A:381:ILE:HD12	1:A:396:VAL:HG23	1.89	0.55
1:A:272:HIS:HD2	1:A:274:GLU:HB2	1.72	0.54
1:A:325:GLN:HE22	1:A:350:ASN:HD21	1.50	0.54
1:A:205:LEU:HB3	1:A:209:GLY:HA2	1.89	0.54
1:B:404:TYR:CE1	1:B:406:LYS:HB2	2.42	0.54
1:A:377:ILE:O	1:A:381:ILE:HG22	2.07	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:B:247:THR:HG21	1:B:297:PRO:HG2	1.90	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:A:384:ASN:CG	1:A:387:LEU:HD12	2.33	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:255:PRO:HA	1:B:259:ASP:OD1	2.08	0.54
1:A:452:MET:HE2	1:A:510:LEU:CD2	2.38	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:HA	1:B:95:GLN:NE2	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:192:ALA:C	1:B:194:ASN:H	2.16	0.53
1:A:406:LYS:HD3	1:A:406:LYS:O	2.08	0.53
1:A:350:ASN:N	1:A:401:VAL:O	2.42	0.53
1:B:468:LYS:HG2	1:B:471:ALA:HB2	1.90	0.53
1:B:122:GLU:CD	1:B:122:GLU:H	2.15	0.53
1:A:287:LEU:HD21	1:A:308:LEU:HD11	1.91	0.53
1:A:322:LYS:HD3	1:A:503:LEU:HD13	1.89	0.53
1:B:16:THR:HG21	1:B:519:GLY:HA3	1.91	0.53
1:B:286:ILE:HD13	1:B:286:ILE:C	2.34	0.53
1:A:58:LEU:HD13	1:A:60:LEU:HD12	1.91	0.52
1:B:244:THR:HG23	1:B:244:THR:O	2.08	0.52
1:B:283:ALA:HA	1:B:286:ILE:HG23	1.91	0.52
1:A:282:ALA:O	1:A:286:ILE:HG13	2.08	0.52
1:A:353:GLY:O	1:A:406:LYS:HG2	2.08	0.52
1:A:150:ASN:HD22	1:A:151:ASN:N	2.06	0.52
1:A:38:THR:HG22	1:A:39:LYS:H	1.74	0.52
1:A:509:ALA:HB1	1:B:531:ARG:NH1	2.17	0.52
1:A:14:VAL:CG2	1:A:518:ILE:HD12	2.39	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:258:ASN:H	1:A:258:ASN:ND2	2.08	0.52
1:A:445:LEU:C	1:A:445:LEU:HD23	2.34	0.52
1:B:193:ALA:C	1:B:195:GLN:H	2.18	0.52
1:A:378:ASP:HA	1:A:381:ILE:HG22	1.91	0.52
1:A:40:THR:HB	1:A:44:ARG:HB3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:158:MET:CE	1:A:176:MET:HE2	2.37	0.51
1:A:418:TYR:OH	1:A:432:ILE:HD12	2.09	0.51
1:B:389:ASN:ND2	1:B:391:LYS:HB3	2.25	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:423:MET:HE3	1:A:424:LEU:CG	2.38	0.51
1:B:299:ASN:H	1:B:299:ASN:ND2	2.07	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:A:273:GLU:HB2	3:A:722:HOH:O	2.11	0.50
1:B:28:THR:O	1:B:55:VAL:HA	2.11	0.50
1:B:158:MET:SD	1:B:176:MET:HE2	2.51	0.50
1:B:523:GLN:HE21	1:B:523:GLN:CA	2.18	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:306:VAL:O	1:B:310:GLU:HG3	2.11	0.50
1:B:467:VAL:O	1:B:468:LYS:HB2	2.12	0.50
1:A:445:LEU:HD21	1:A:487:TRP:HB3	1.94	0.50
1:B:134:LEU:HB3	1:B:135:PRO:HD2	1.93	0.50
1:B:192:ALA:C	1:B:194:ASN:N	2.68	0.50
1:A:469:GLU:HG2	1:A:470:ASP:N	2.26	0.50
1:B:122:GLU:CD	1:B:122:GLU:N	2.70	0.50
1:A:158:MET:HE2	1:A:176:MET:CG	2.39	0.50
1:A:258:ASN:H	1:A:258:ASN:HD22	1.58	0.50
1:B:95:GLN:HE21	1:B:95:GLN:CA	2.25	0.50
1:B:97:LYS:NZ	1:B:97:LYS:CB	2.75	0.50
1:A:88:ASN:HD21	1:A:104:ASN:HA	1.76	0.49
1:A:100:VAL:O	1:A:100:VAL:HG13	2.11	0.49
1:A:426:GLY:HA3	1:B:440:LEU:CD1	2.42	0.49
1:B:130:PHE:CE2	1:B:136:MET:HE1	2.47	0.49
1:B:285:SER:O	1:B:290:VAL:HG22	2.12	0.49
1:A:435:VAL:HG13	1:B:429:ARG:HB3	1.94	0.49
1:B:332:LEU:CD1	1:B:432:ILE:HD11	2.17	0.49
1:B:389:ASN:HD21	1:B:391:LYS:HB3	1.77	0.49
1:A:426:GLY:HA3	1:B:440:LEU:HD13	1.94	0.49
1:B:258:ASN:HD22	1:B:258:ASN:N	2.03	0.49
1:B:413:VAL:HA	1:B:434:ASN:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:LEU:HD22	1:A:305:LEU:HA	1.93	0.49
1:B:337:VAL:HG21	1:B:380:ILE:HG22	1.95	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
1:B:405:MET:O	1:B:408:VAL:HG23	2.12	0.49
1:A:153:ASP:OD2	1:A:153:ASP:C	2.54	0.49
1:B:252:GLU:O	1:B:274:GLU:OE1	2.29	0.49
1:B:381:ILE:HD11	1:B:396:VAL:CG2	2.43	0.49
1:B:337:VAL:HG21	1:B:380:ILE:CG2	2.43	0.49
1:B:454:GLU:C	1:B:454:GLU:CD	2.80	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:335:PHE:HZ	1:B:327:LYS:HE2	1.77	0.48
1:A:343:PRO:HA	1:A:420:SER:HA	1.94	0.48
1:A:351:HIS:HA	1:A:403:LYS:O	2.12	0.48
1:B:238:LYS:HZ2	1:B:457:THR:HG21	1.78	0.48
1:B:218:THR:HG23	3:B:699:HOH:O	2.12	0.48
1:B:249:ARG:HG2	1:B:249:ARG:NH2	2.27	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:A:143:VAL:HG21	1:A:236:LEU:HD21	1.96	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:B:523:GLN:HA	1:B:523:GLN:NE2	2.24	0.48
1:A:104:ASN:HD22	1:A:104:ASN:C	2.21	0.48
1:A:200:ASN:N	1:A:200:ASN:ND2	2.62	0.48
1:A:234:ASN:HB2	1:A:236:LEU:HG	1.95	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: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:A:32:TYR:OH	1:B:529:GLU:OE2	2.28	0.48
1:A:50:THR:HG22	3:A:677:HOH:O	2.14	0.48
1:B:195:GLN:HE22	1:B:198:ARG:HD2	1.79	0.48
1:A:56:PHE:CE1	1:A:461:TYR:HB3	2.49	0.47
1:A:481:LEU:HD13	1:A:484:LEU:HD21	1.95	0.47
1:B:131:ASN:HD22	1:B:131:ASN:N	2.12	0.47
1:B:82:VAL:HG21	1:B:154:LEU:CD2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:CYS:HB3	1:A:28:THR:HA	1.96	0.47
1:A:412:LYS:HB3	1:A:412:LYS:NZ	2.25	0.47
1:A:511:GLU:O	1:A:515:ARG:HG3	2.14	0.47
1:A:158:MET:HG3	1:A:176:MET:HE2	1.97	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:B:526:LEU:O	1:B:527:ARG:C	2.57	0.47
1:A:247:THR:CG2	1:A:297:PRO:HG2	2.45	0.47
1:A:406:LYS:HB3	1:A:407:PRO:HD3	1.97	0.47
1:A:533:LEU:HD22	1:A:533:LEU:H	1.79	0.47
1:B:229:ASN:O	1:B:233:GLU:HB3	2.14	0.47
1:A:155:TYR:O	1:A:159:GLN:HG3	2.14	0.47
1:A:476:ASN:O	1:A:477:PHE:C	2.58	0.47
1:B:208:LYS:CB	1:B:210:ASN:HD22	2.28	0.47
1:A:67:GLY:HA3	1:A:239:VAL:HG22	1.96	0.47
1:A:88:ASN:HD21	1:A:104:ASN:CA	2.28	0.47
1:A:154:LEU:HD22	1:A:179:VAL:CG1	2.44	0.47
1:B:453:THR:HG23	1:B:481:LEU:CD1	2.44	0.47
1:A:220:LEU:O	1:A:224:ARG:HG3	2.15	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:412:LYS:NZ	1:A:414:ALA:HB2	2.29	0.47
1:B:293:ILE:HG23	1:B:317:ALA:CB	2.45	0.47
1:B:26:LEU:HD12	1:B:27:LEU:H	1.79	0.46
1:A:497:PHE:CE2	1:B:530:GLU:HB2	2.50	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:B:524:ASN:O	1:B:525:GLU:HB2	2.14	0.46
1:A:516:LEU:C	1:A:516:LEU:CD2	2.88	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: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:B:353:GLY:HA3	1:B:409:GLY:O	2.16	0.46
1:A:200:ASN:HD22	1:A:200:ASN:C	2.23	0.46
1:A:378:ASP:HA	1:A:381:ILE:CG2	2.46	0.46
1:A:158:MET:HE1	1:A:172:LEU:HB2	1.97	0.46
1:B:258:ASN:N	1:B:258:ASN:ND2	2.63	0.46
1:A:256:GLY:HA2	1:A:263:ASN:ND2	2.31	0.46
1:B:380:ILE:HG22	1:B:380:ILE:O	2.16	0.46
1:A:286:ILE:HA	1:A:314:THR:HG21	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ASN:HB3	1:B:439:SER:OG	2.16	0.45
1:B:134:LEU:HB3	1:B:135:PRO:CD	2.47	0.45
1:B:131:ASN:N	1:B:131:ASN:ND2	2.64	0.45
1:A:14:VAL:HG22	1:B:47:VAL:CG1	2.47	0.45
1:A:55:VAL:HG23	1:A:464:VAL:HG21	1.99	0.45
1:B:97:LYS:HB2	1:B:97:LYS:HZ2	1.80	0.45
1:B:465:ASP:OD2	1:B:467:VAL:HB	2.16	0.45
1:A:223:ILE:O	1:A:226:ASP:HB2	2.17	0.45
1:A:258:ASN:ND2	1:A:258:ASN:N	2.62	0.45
1:A:335:PHE:CZ	1:B:327:LYS:HE2	2.51	0.45
1:B:186:TYR:CZ	1:B:188:PRO:HB3	2.52	0.45
1:B:477:PHE:O	1:B:478:TYR:C	2.59	0.45
1:B:389:ASN:OD1	1:B:392:LEU:HD13	2.17	0.45
1:B:529:GLU:CD	1:B:529:GLU:H	2.25	0.45
1:A:44:ARG:HH21	1:B:13:LYS:CG	2.27	0.45
1:A:104:ASN:ND2	1:A:104:ASN:C	2.74	0.45
1:A:337:VAL:HG21	1:A:380:ILE:HG22	1.98	0.45
1:B:200:ASN:HD22	1:B:200:ASN:H	1.65	0.45
1:B:465:ASP:CG	1:B:466:PRO:HD2	2.42	0.45
1:A:122:GLU:OE2	1:A:122:GLU:N	2.50	0.45
1:A:352:LEU:N	1:A:352:LEU:CD2	2.79	0.45
1:A:183:PRO:HB2	1:A:203:ILE:HG23	2.00	0.44
1:A:294:ASN:ND2	1:A:294:ASN:C	2.72	0.44
1:A:325:GLN:HG3	3:A:652:HOH:O	2.16	0.44
1:B:150:ASN:ND2	1:B:160:ARG:NH1	2.64	0.44
1:A:115:LEU:HD22	1:A:511:GLU:HB3	2.00	0.44
1:A:286:ILE:HG23	1:A:314:THR:HB	1.98	0.44
1:A:378:ASP:CA	1:A:381:ILE:HG22	2.47	0.44
1:A:351:HIS:O	1:A:351:HIS:ND1	2.51	0.44
1:B:285:SER:HB3	1:B:290:VAL:HG23	1.99	0.44
1:B:302:VAL:HG23	1:B:305:LEU:H	1.82	0.44
1:A:301:PHE:HA	1:A:305:LEU:HD23	1.98	0.44
1:B:26:LEU:HD12	1:B:27:LEU:N	2.33	0.44
1:B:272:HIS:ND1	1:B:274:GLU:N	2.59	0.44
1:B:352:LEU:N	1:B:352:LEU:CD2	2.81	0.44
1:A:465:ASP:OD1	1:A:467:VAL:HG22	2.18	0.44
1:A:477:PHE:O	1:A:478:TYR:C	2.60	0.44
1:B:293:ILE:CA	1:B:317:ALA:HB2	2.35	0.44
1:A:57:LYS:HB3	1:A:460:SER:HB3	1.98	0.44
1:A:169:GLN:HE21	1:A:169:GLN:HB3	1.61	0.44
1:A:500:VAL:HG21	1:B:527:ARG:NH2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ASN:HD22	2:B:660:NAD:H8A	1.81	0.44
1:B:438:ASP:HB3	2:B:660:NAD:O7N	2.18	0.44
1:A:203:ILE:C	1:A:203:ILE:HD12	2.43	0.44
1:B:174:ALA:HB3	3:B:669:HOH:O	2.18	0.44
1:B:293:ILE:HA	1:B:317:ALA:HB3	1.93	0.44
1:A:52:GLN:OE1	1:A:53:ASP:O	2.36	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:B:350:ASN:O	1:B:351:HIS:HB3	2.18	0.44
1:A:109:MET:HE2	1:A:507:ARG:HE	1.83	0.43
1:A:302:VAL:HB	1:A:303:PRO:HD2	2.00	0.43
1:A:530:GLU:HG3	1:B:497:PHE:HD2	1.82	0.43
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: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:A:440:LEU:HD12	1:A:440:LEU:HA	1.83	0.43
1:A:447:ILE:O	1:A:451:VAL:HG23	2.19	0.43
1:A:520:LEU:HA	1:A:520:LEU:HD23	1.77	0.43
1:A:525:GLU:HA	1:A:525:GLU:OE2	2.18	0.43
1:A:532:LEU:HG	1:B:461:TYR:CZ	2.54	0.43
1:B:116:LYS:C	1:B:116:LYS:HD2	2.44	0.43
1:B:254:SER:O	1:B:257:VAL:HG13	2.18	0.43
1:A:38:THR:O	1:A:45:PHE:HA	2.19	0.43
1:A:47:VAL:HG11	1:B:520:LEU:HD11	2.01	0.43
1:A:125:ASP:OD1	1:A:527:ARG:NH2	2.51	0.43
1:A:134:LEU:HD21	1:A:518:ILE:CG2	2.48	0.43
1:A:190:PHE:HB3	1:A:248:GLU:HB3	2.01	0.43
1:A:200:ASN:ND2	1:A:200:ASN:C	2.76	0.43
1:B:260:THR:OG1	1:B:263:ASN:ND2	2.52	0.43
1:A:234:ASN:CB	1:A:236:LEU:HG	2.49	0.43
1:A:354:ASN:HD21	1:A:356:ASP:HB2	1.83	0.43
1:B:115:LEU:O	1:B:127:TYR:HA	2.19	0.43
1:A:51:VAL:HG12	1:A:52:GLN:N	2.34	0.42
1:B:276:ALA:O	1:B:279:THR:HB	2.19	0.42
1:B:350:ASN:HD21	1:B:402:ILE:CG1	2.32	0.42
1:A:14:VAL:HG22	1:B:47:VAL:HG11	2.00	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:B:381:ILE:HG23	1:B:388:TYR:CG	2.54	0.42
1:A:349:TYR:HD1	1:A:401:VAL:HB	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:MET:O	1:A:406:LYS:C	2.63	0.42
1:A:483:PHE:CE1	1:B:531:ARG:HD3	2.55	0.42
1:B:163:VAL:HG23	1:B:164:LEU:HD13	2.01	0.42
1:B:350:ASN:ND2	1:B:402:ILE:HG13	2.35	0.42
1:B:501:ASN:HA	1:B:506:GLN:OE1	2.19	0.42
1:A:10:THR:HG21	1:A:129:PRO:HG2	2.00	0.42
1:A:242:LEU:HD12	1:A:293:ILE:HD12	2.01	0.42
1:A:247:THR:HG23	1:A:298:GLN:HE21	1.84	0.42
1:A:250:TYR:CD2	1:A:299:ASN:OD1	2.73	0.42
1:B:71:ILE:HG13	1:B:147:TRP:CE3	2.54	0.42
1:B:473:LYS:CD	1:B:474:PHE:H	2.30	0.42
1:A:284:ALA:O	1:A:285:SER:C	2.61	0.42
1:B:195:GLN:NE2	1:B:195:GLN:HA	2.35	0.42
1:B:207:GLU:C	1:B:209:GLY:N	2.77	0.42
1:B:248:GLU:OE1	1:B:277:PRO:HD2	2.20	0.42
1:B:258:ASN:H	1:B:258:ASN:ND2	2.13	0.42
1:A:122:GLU:N	1:A:122:GLU:CD	2.76	0.42
1:A:299:ASN:O	1:A:299:ASN:CG	2.60	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:A:258:ASN:HD22	1:A:258:ASN:C	2.27	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: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: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:B:56:PHE:HA	1:B:461:TYR:HA	2.02	0.42
1:B:469:GLU:O	1:B:470:ASP:C	2.63	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:NE2	1:B:95:GLN:CA	2.83	0.41
1:B:208:LYS:HB2	1:B:210:ASN:HD22	1.85	0.41
1:B:245:ALA:O	1:B:246:ASN:C	2.61	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:A:73:LEU:HD23	1:A:181:PRO:HB3	2.01	0.41
1:A:412:LYS:HG3	1:A:438:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:ILE:HG21	1:B:243:TRP:CZ3	2.56	0.41
1:B:468:LYS:HZ3	1:B:471:ALA:CB	2.34	0.41
1:A:56:PHE:CD1	1:A:461:TYR:HB3	2.55	0.41
1:A:150:ASN:ND2	1:A:151:ASN:N	2.68	0.41
1:A:325:GLN:HE21	1:A:325:GLN:HB3	1.68	0.41
1:B:231:LYS:HG2	1:B:239:VAL:HG21	2.03	0.41
1:B:268:ILE:HG12	1:B:275:ILE:HD12	2.03	0.41
1:A:10:THR:HG23	1:A:132:SER:CB	2.50	0.41
1:A:67:GLY:CA	1:A:236:LEU:HD13	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: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:B:70:LEU:HD12	1:B:242:LEU:HB3	2.02	0.41
1:A:175:LYS:HA	1:A:175:LYS:HD3	1.83	0.41
1:A:320:ASP:HB3	1:A:489:LYS:HB3	2.01	0.41
1:A:406:LYS:N	1:A:407:PRO:CD	2.83	0.41
1:B:130:PHE:HE2	1:B:136:MET:HE1	1.83	0.41
1:A:109:MET:HE2	1:A:507:ARG:NE	2.36	0.41
1:A:215:GLY:HA2	3:A:660:HOH:O	2.21	0.41
1:A:328:LEU:HD11	1:B:332:LEU:HG	2.02	0.41
1:B:195:GLN:NE2	1:B:198:ARG:HD2	2.36	0.41
1:B:254:SER:O	1:B:258:ASN:ND2	2.54	0.41
1:B:424:LEU:HA	3:B:702:HOH:O	2.19	0.41
1:B:486:TYR:CE1	1:B:507:ARG:HB2	2.56	0.41
1:A:21:TYR:CZ	1:A:26:LEU:HD13	2.56	0.41
1:A:174:ALA:HB3	3:A:709:HOH:O	2.20	0.41
1:A:247:THR:HG23	1:A:298:GLN:NE2	2.36	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:381:ILE:HG13	1:A:388:TYR:CD2	2.55	0.41
1:A:507:ARG:O	1:A:510:LEU:HB2	2.21	0.41
1:A:532:LEU:HD23	1:B:54:TYR:CE2	2.55	0.41
1:B:119:ILE:HA	1:B:124:ASN:O	2.21	0.41
1:B:323:SER:O	1:B:327:LYS:HG3	2.21	0.41
1:B:349:TYR:CD1	1:B:349:TYR:N	2.88	0.41
1:A:10:THR:HG23	1:A:132:SER:HB2	2.02	0.40
1:A:195:GLN:NE2	1:A:355:ASN:HD21	2.15	0.40
1:A:228:GLN:HG2	1:A:288:GLU:OE1	2.21	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:LEU:CB	1:A:60:LEU:HD11	2.52	0.40
1:A:71:ILE:HG21	1:A:243:TRP:CZ3	2.56	0.40
1:A:208:LYS:O	1:A:209:GLY:C	2.65	0.40
1:B:245:ALA:HA	2:B:660:NAD:H52A	2.02	0.40
1:B:483:PHE:HD2	1:B:483:PHE:HA	1.81	0.40
1:A:418:TYR:CD1	1:A:418:TYR:N	2.90	0.40
1:B:322:LYS:HD2	1:B:489:LYS:HA	2.03	0.40
1:B:484:LEU:HA	1:B:487:TRP:CZ3	2.57	0.40
1:B:507:ARG:NH2	1:B:507:ARG:HG3	2.30	0.40
1:A:173:LYS:O	1:A:174:ALA:C	2.64	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

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	502/533 (94%)	431 (86%)	57 (11%)	14 (3%)	4	3
1	B	494/533 (93%)	429 (87%)	52 (10%)	13 (3%)	4	4
All	All	996/1066 (93%)	860 (86%)	109 (11%)	27 (3%)	4	4

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

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Mol	Chain	Res	Type
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
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%)	412 (92%)	35 (8%)	11	20
1	B	439/471 (93%)	394 (90%)	45 (10%)	7	11
All	All	886/942 (94%)	806 (91%)	80 (9%)	9	15

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU

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Mol	Chain	Res	Type
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	154	LEU
1	A	189	ASP
1	A	200	ASN
1	A	207	GLU
1	A	211	VAL
1	A	258	ASN
1	A	260	THR
1	A	262	GLU
1	A	290	VAL
1	A	298	GLN
1	A	300	THR
1	A	320	ASP
1	A	332	LEU
1	A	336	LEU
1	A	352	LEU
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	438	ASP
1	A	473	LYS
1	A	516	LEU
1	A	532	LEU
1	B	14	VAL
1	B	70	LEU
1	B	73	LEU
1	B	92	VAL
1	B	95	GLN
1	B	97	LYS
1	B	111	GLN
1	B	116	LYS

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Mol	Chain	Res	Type
1	B	122	GLU
1	B	131	ASN
1	B	151	ASN
1	B	163	VAL
1	B	164	LEU
1	B	168	LEU
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	257	VAL
1	B	258	ASN
1	B	262	GLU
1	B	286	ILE
1	B	299	ASN
1	B	308	LEU
1	B	349	TYR
1	B	352	LEU
1	B	387	LEU
1	B	391	LYS
1	B	394	LYS
1	B	408	VAL
1	B	441	LEU
1	B	457	THR
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	508	THR
1	B	517	LEU
1	B	523	GLN
1	B	529	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (59) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	ASN

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Mol	Chain	Res	Type
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	194	ASN
1	A	195	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
1	A	325	GLN
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	102	GLN
1	B	111	GLN
1	B	131	ASN
1	B	150	ASN
1	B	151	ASN
1	B	159	GLN
1	B	195	GLN
1	B	200	ASN
1	B	201	ASN
1	B	210	ASN
1	B	219	HIS

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Mol	Chain	Res	Type
1	B	221	GLN
1	B	228	GLN
1	B	229	ASN
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	334	GLN
1	B	350	ASN
1	B	476	ASN
1	B	498	HIS
1	B	504	ASN
1	B	506	GLN
1	B	523	GLN

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 oligosaccharides 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
2	NAD	B	660	-	46,48,48	2.32	13 (28%)	64,73,73	1.09	7 (10%)
2	NAD	A	650	-	46,48,48	2.12	9 (19%)	64,73,73	1.04	6 (9%)

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	B	660	-	-	7/30/62/62	0/5/5/5
2	NAD	A	650	-	-	3/30/62/62	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	660	NAD	C2N-N1N	8.46	1.44	1.35
2	B	660	NAD	C4N-C3N	7.13	1.50	1.39
2	A	650	NAD	C4N-C3N	6.87	1.49	1.39
2	A	650	NAD	C2N-N1N	6.65	1.42	1.35
2	B	660	NAD	C6N-N1N	4.68	1.46	1.35
2	A	650	NAD	C6N-N1N	4.58	1.45	1.35
2	A	650	NAD	C5N-C4N	4.39	1.46	1.38
2	B	660	NAD	C2A-N1A	4.17	1.41	1.33
2	B	660	NAD	C5N-C4N	4.05	1.45	1.38
2	A	650	NAD	C2A-N1A	4.04	1.41	1.33
2	B	660	NAD	PA-O3	-3.29	1.55	1.59
2	A	650	NAD	PN-O3	2.73	1.62	1.59
2	B	660	NAD	O4B-C1B	-2.68	1.35	1.42
2	B	660	NAD	C3N-C7N	2.61	1.54	1.50
2	B	660	NAD	C8A-N7A	2.59	1.36	1.31
2	B	660	NAD	C5A-C4A	2.37	1.43	1.39
2	B	660	NAD	C2N-C3N	2.30	1.42	1.39
2	A	650	NAD	C5A-C6A	2.29	1.47	1.41
2	A	650	NAD	C8A-N7A	2.21	1.35	1.31
2	B	660	NAD	C5A-C6A	2.19	1.47	1.41
2	B	660	NAD	C4A-N3A	2.15	1.38	1.34
2	A	650	NAD	C3N-C7N	2.02	1.53	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	660	NAD	C3N-C7N-N7N	3.12	121.58	117.74
2	A	650	NAD	O2A-PA-O3	2.97	115.31	107.27
2	A	650	NAD	C3N-C7N-N7N	2.89	121.30	117.74
2	B	660	NAD	O2A-PA-O3	2.72	114.64	107.27
2	B	660	NAD	C6N-N1N-C2N	-2.64	119.63	121.88
2	B	660	NAD	O7N-C7N-N7N	-2.48	119.03	122.62
2	A	650	NAD	O4D-C4D-C3D	2.39	109.89	105.15
2	A	650	NAD	O7N-C7N-N7N	-2.28	119.32	122.62
2	B	660	NAD	O3-PN-O1N	-2.22	104.03	110.70
2	B	660	NAD	O4B-C4B-C5B	-2.13	102.51	109.33
2	B	660	NAD	C4D-O4D-C1D	-2.13	107.97	109.92
2	A	650	NAD	O3-PN-O1N	-2.05	104.54	110.70
2	A	650	NAD	C4A-N9A-C1B	-2.02	121.90	126.63

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	650	NAD	PN-O3-PA-O5B
2	B	660	NAD	C5B-O5B-PA-O1A
2	B	660	NAD	C5B-O5B-PA-O2A
2	B	660	NAD	C5B-O5B-PA-O3
2	A	650	NAD	PA-O3-PN-O1N
2	B	660	NAD	PA-O3-PN-O2N
2	B	660	NAD	O4B-C4B-C5B-O5B
2	A	650	NAD	PA-O3-PN-O2N
2	B	660	NAD	PA-O3-PN-O1N
2	B	660	NAD	C3B-C4B-C5B-O5B

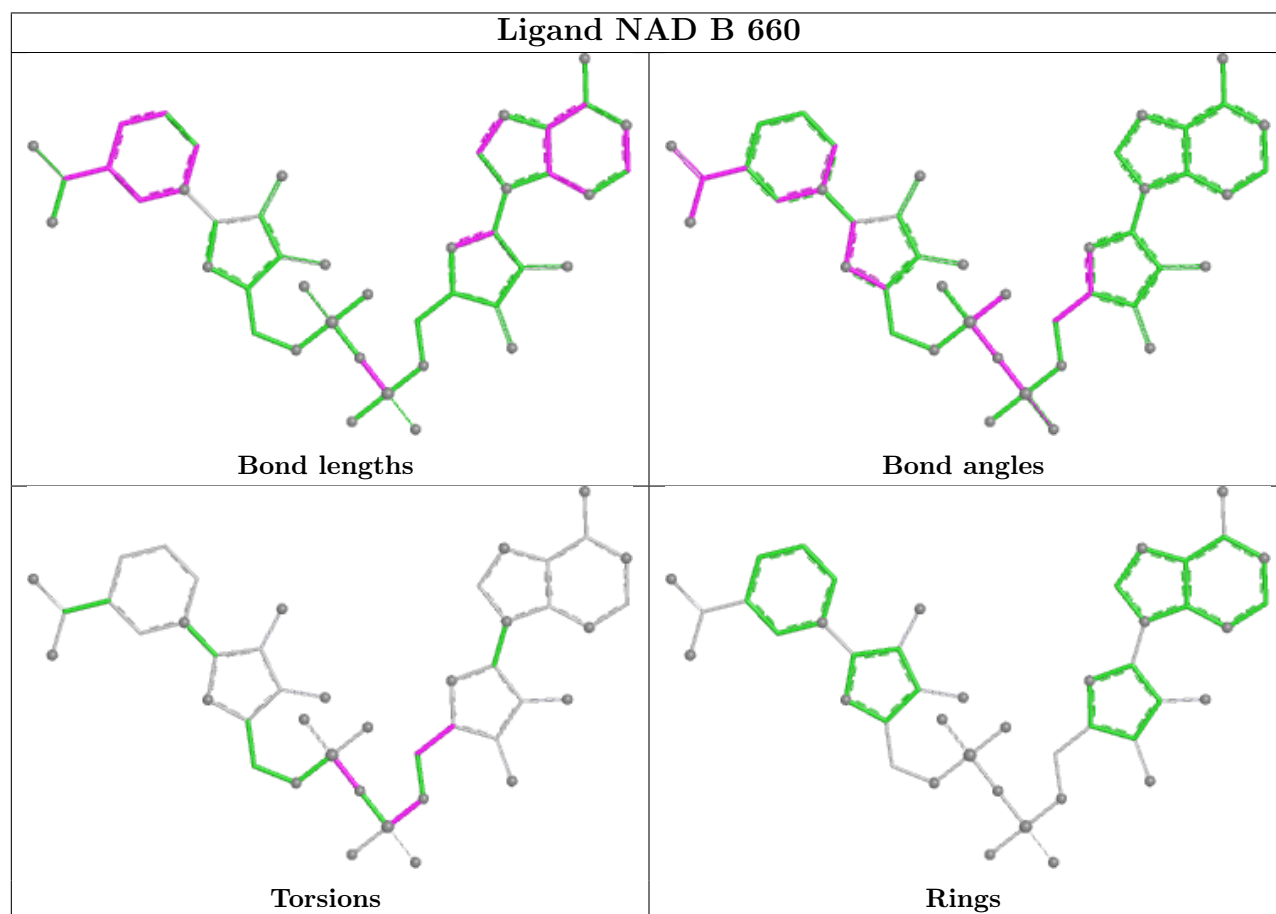
There are no ring outliers.

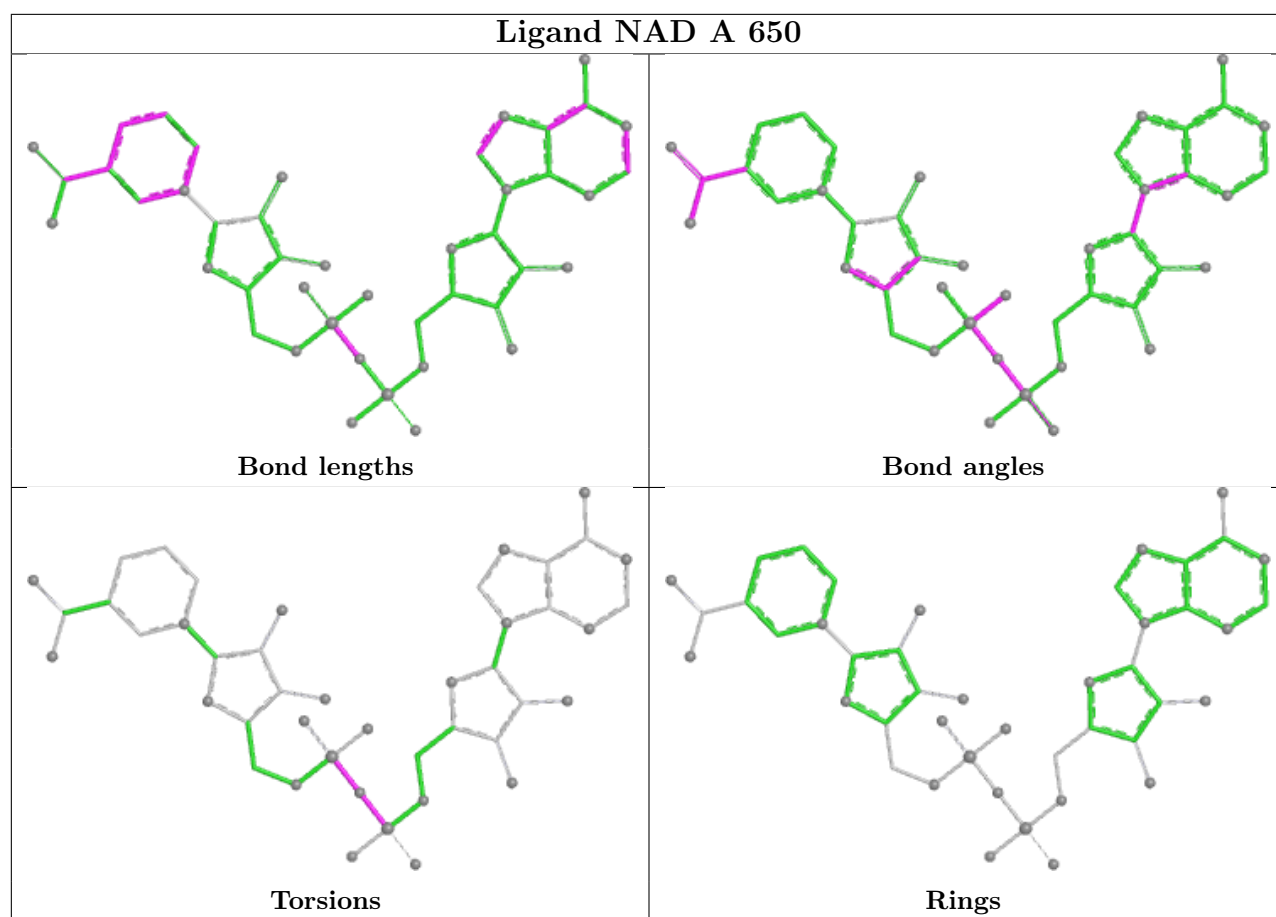
2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	660	NAD	5	0
2	A	650	NAD	2	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 ⓘ

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%)	0.03	10 (1%) 65 60	26, 53, 91, 91	0
1	B	498/533 (93%)	0.17	23 (4%) 37 33	27, 56, 91, 91	0
All	All	1004/1066 (94%)	0.10	33 (3%) 49 45	26, 54, 91, 91	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	467	VAL	4.1
1	B	349	TYR	3.8
1	A	330	SER	3.5
1	B	352	LEU	3.4
1	B	469	GLU	3.2
1	A	396	VAL	3.0
1	A	400	ILE	3.0
1	B	393	GLY	3.0
1	B	197	GLU	3.0
1	B	193	ALA	3.0
1	A	43	GLY	3.0
1	B	468	LYS	2.7
1	A	404	TYR	2.7
1	B	402	ILE	2.7
1	B	209	GLY	2.6
1	B	317	ALA	2.5
1	B	380	ILE	2.4
1	B	404	TYR	2.4
1	B	348	SER	2.4
1	B	409	GLY	2.3
1	A	412	LYS	2.3
1	B	206	ASP	2.3
1	B	347	ALA	2.2
1	A	351	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	392	LEU	2.2
1	A	259	ASP	2.1
1	A	391	LYS	2.1
1	B	410	ASP	2.1
1	B	391	LYS	2.0
1	B	408	VAL	2.0
1	B	192	ALA	2.0
1	B	397	ASP	2.0
1	A	460	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

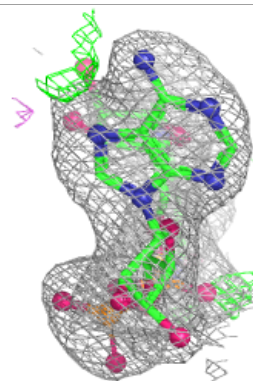
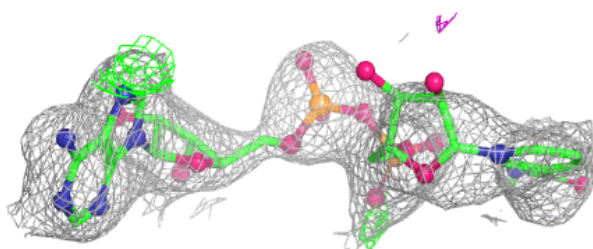
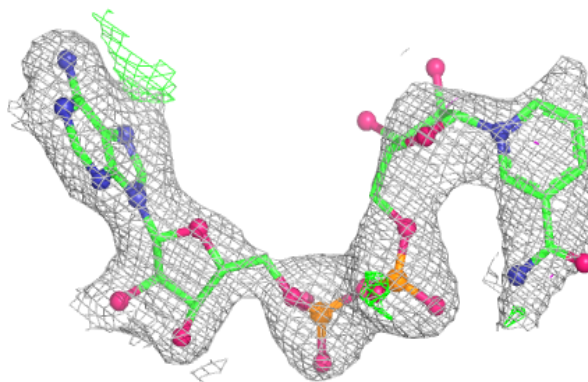
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAD	A	650	44/44	0.95	0.09	36,57,84,86	0
2	NAD	B	660	44/44	0.95	0.09	43,56,79,81	0

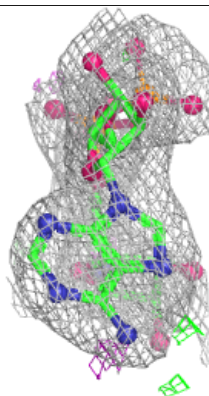
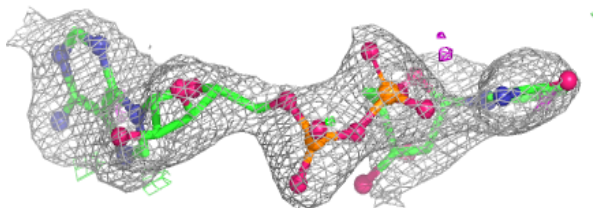
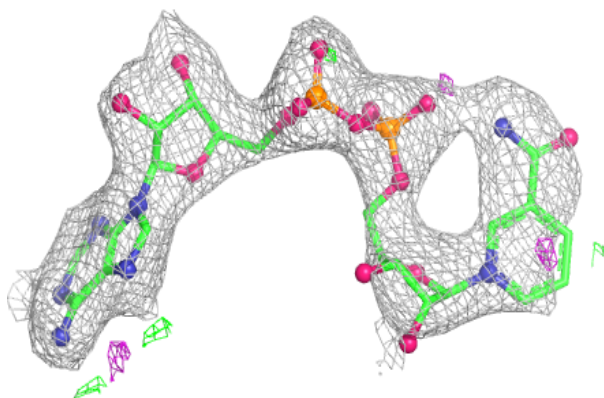
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around NAD A 650:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around NAD B 660:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.