



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:50 AM UTC

PDB ID : 1AG8 / pdb_00001ag8
Title : ALDEHYDE DEHYDROGENASE FROM BOVINE MITOCHONDRIA
Authors : Steinmetz, C.G.; Hurley, T.D.
Deposited on : 1997-04-03
Resolution : 2.65 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
Xtriage (Phenix) : 2.0
EDS : 3.0
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

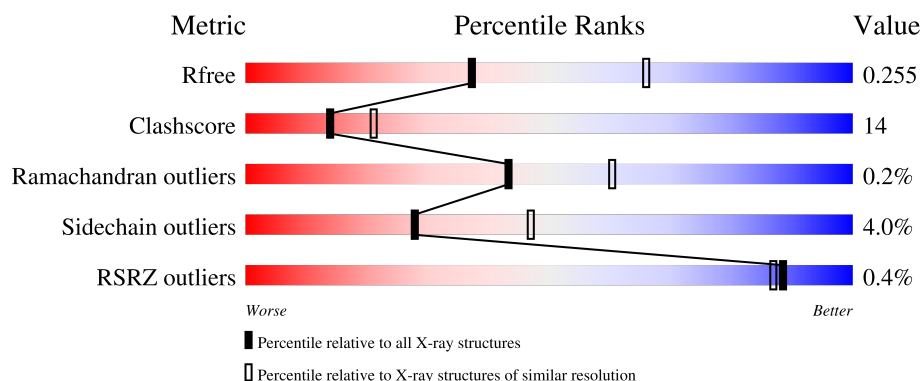
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

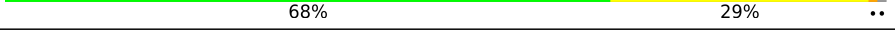
The reported resolution of this entry is 2.65 Å.

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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	499	 68% 28% ..
1	B	499	 70% 25% ..
1	C	499	 67% 30% ..
1	D	499	 68% 29% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15336 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 ALDEHYDE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	0	0	0
			3799	2418	650	714	17			
1	B	493	Total	C	N	O	S	0	0	0
			3799	2418	650	714	17			
1	C	493	Total	C	N	O	S	0	0	0
			3799	2418	650	714	17			
1	D	493	Total	C	N	O	S	0	0	0
			3799	2418	650	714	17			

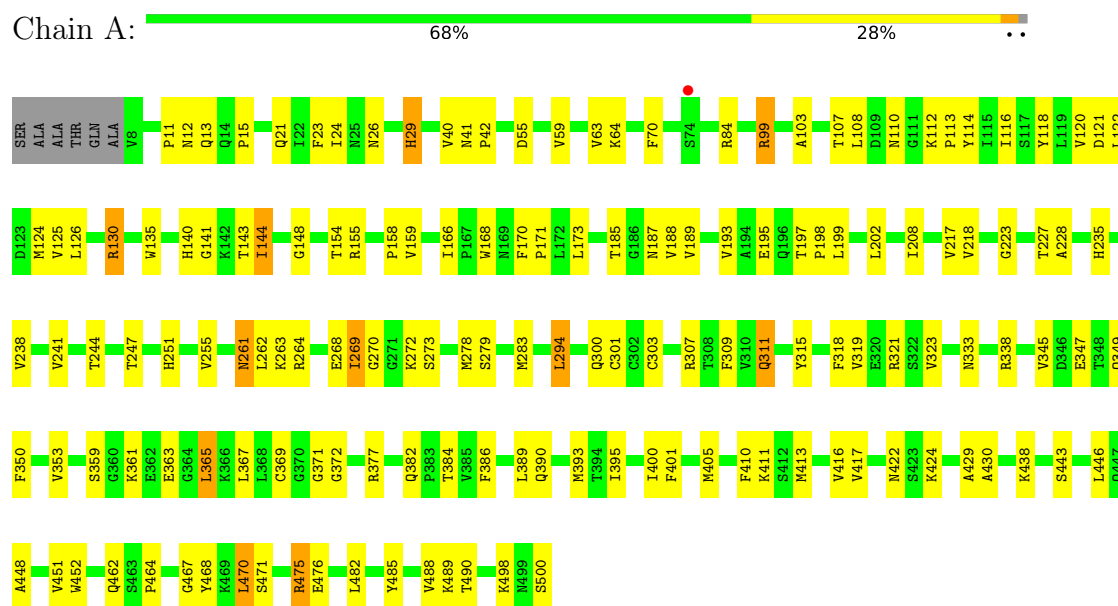
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	34	Total	O	0	0
			34	34		
2	B	37	Total	O	0	0
			37	37		
2	C	34	Total	O	0	0
			34	34		
2	D	35	Total	O	0	0
			35	35		

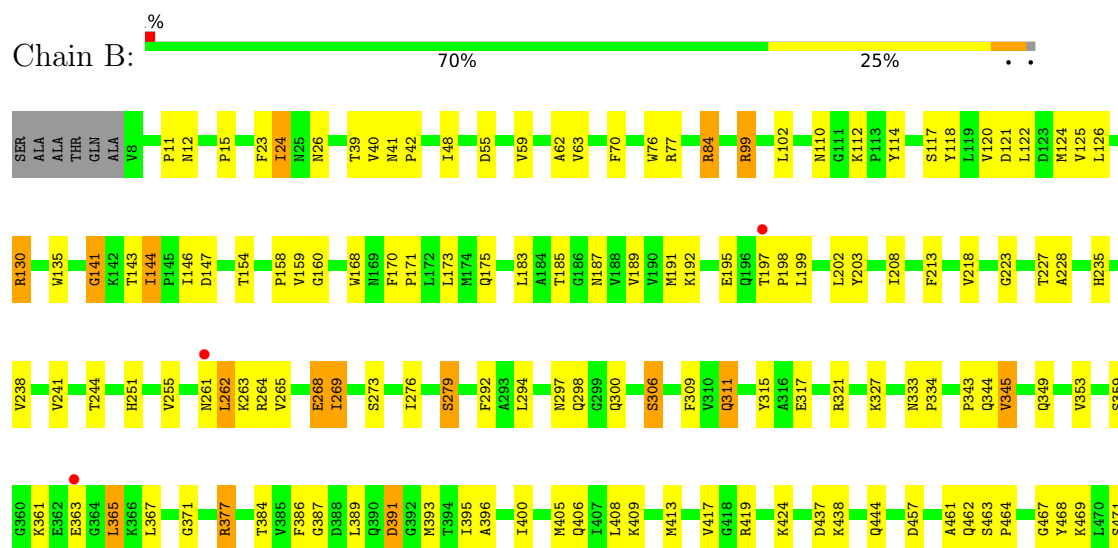
3 Residue-property plots

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: ALDEHYDE DEHYDROGENASE



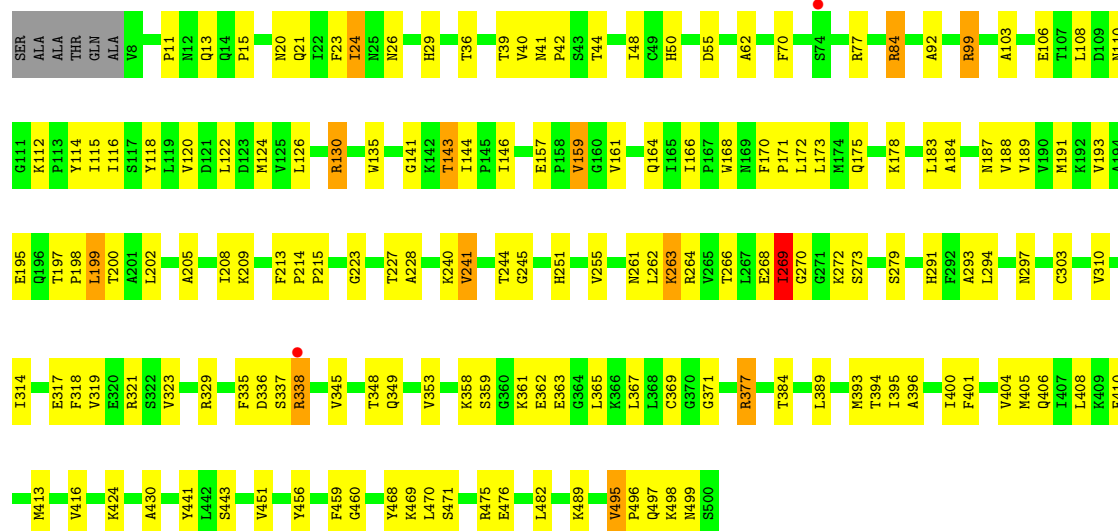
• Molecule 1: ALDEHYDE DEHYDROGENASE





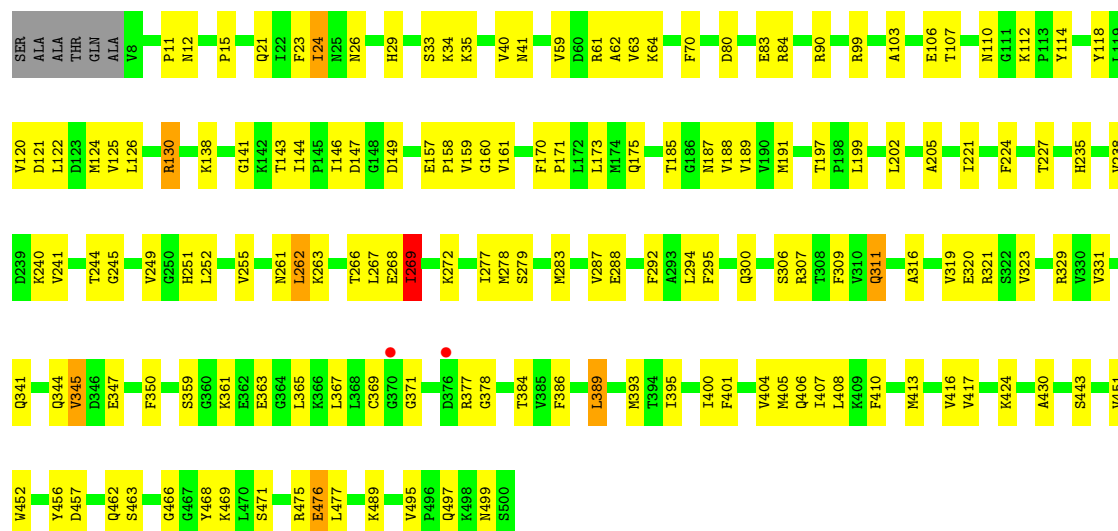
• Molecule 1: ALDEHYDE DEHYDROGENASE

Chain C: 67% 30% ..



• Molecule 1: ALDEHYDE DEHYDROGENASE

Chain D: 68% 29% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	122.60Å 198.40Å 91.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.65 8.00 – 2.65	Depositor EDS
% Data completeness (in resolution range)	83.1 (8.00-2.65) 79.9 (8.00-2.65)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.54 (at 2.65Å)	Xtriage
Refinement program	XTALVIEW, X-PLOR 3.1	Depositor
R, R_{free}	0.217 , 0.282 0.200 , 0.255	Depositor DCC
R_{free} test set	3941 reflections (7.10%)	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	15336	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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

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	0.48	0/3884	0.96	13/5268 (0.2%)
1	B	0.48	0/3884	0.97	14/5268 (0.3%)
1	C	0.44	0/3884	0.94	11/5268 (0.2%)
1	D	0.47	0/3884	0.96	11/5268 (0.2%)
All	All	0.47	0/15536	0.96	49/21072 (0.2%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	144	ILE	N-CA-C	8.72	117.40	107.89
1	D	144	ILE	N-CA-C	8.44	117.09	107.89
1	B	199	LEU	N-CA-C	8.07	119.70	111.07
1	C	476	GLU	N-CA-C	7.88	123.08	113.38
1	A	166	ILE	N-CA-C	7.87	116.58	107.84
1	C	141	GLY	N-CA-C	-7.50	101.87	112.60
1	A	199	LEU	N-CA-C	7.46	120.13	111.02
1	A	144	ILE	N-CA-C	7.46	116.02	107.89
1	B	476	GLU	N-CA-C	7.43	122.52	113.38
1	C	166	ILE	N-CA-C	7.37	116.02	107.84
1	D	141	GLY	N-CA-C	-7.20	102.31	112.60
1	C	189	VAL	N-CA-C	6.90	118.41	108.48
1	B	189	VAL	N-CA-C	6.89	118.40	108.48
1	C	273	SER	N-CA-C	6.87	118.02	109.57
1	C	199	LEU	N-CA-C	6.84	119.37	111.02
1	B	141	GLY	N-CA-C	-6.80	102.88	112.60
1	D	199	LEU	N-CA-C	6.76	119.27	111.02
1	C	144	ILE	N-CA-C	6.75	116.01	108.05
1	A	476	GLU	N-CA-C	6.49	122.16	113.72
1	B	273	SER	N-CA-C	6.44	118.76	109.48
1	A	141	GLY	N-CA-C	-6.41	103.43	112.60
1	D	476	GLU	N-CA-C	6.37	121.33	113.50
1	A	333	ASN	CA-C-N	6.31	126.79	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	ASN	C-N-CA	6.31	126.79	119.47
1	D	345	VAL	N-CA-C	6.11	117.58	110.62
1	C	84	ARG	N-CA-C	-6.04	104.70	111.28
1	D	189	VAL	N-CA-C	5.98	117.37	108.46
1	A	189	VAL	N-CA-C	5.92	117.01	108.48
1	B	143	THR	N-CA-C	-5.91	98.89	108.52
1	A	143	THR	N-CA-C	-5.83	99.02	108.52
1	D	84	ARG	N-CA-C	-5.52	105.27	111.28
1	A	84	ARG	N-CA-C	-5.48	105.31	111.28
1	A	273	SER	N-CA-C	5.48	117.37	109.48
1	B	345	VAL	N-CA-C	5.44	116.82	110.62
1	D	261	ASN	N-CA-C	5.43	119.62	112.89
1	B	84	ARG	N-CA-C	-5.36	105.43	111.28
1	D	344	GLN	N-CA-C	-5.29	102.64	110.48
1	B	344	GLN	N-CA-C	-5.26	102.69	110.48
1	B	391	ASP	N-CA-C	5.22	117.65	111.33
1	C	261	ASN	N-CA-C	5.21	119.35	112.89
1	D	143	THR	N-CA-C	-5.14	100.14	108.52
1	A	140	HIS	N-CA-C	5.12	117.13	110.43
1	B	437	ASP	N-CA-C	5.08	116.82	111.28
1	D	269	ILE	N-CA-C	5.07	119.88	109.34
1	C	143	THR	N-CA-C	-5.06	100.49	108.73
1	B	276	ILE	N-CA-C	5.05	115.60	107.73
1	A	261	ASN	N-CA-C	5.02	119.12	112.89
1	B	261	ASN	N-CA-C	5.02	119.11	112.89
1	C	269	ILE	N-CA-C	5.01	119.76	109.34

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	3799	0	3755	105	0
1	B	3799	0	3755	115	0
1	C	3799	0	3755	113	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3799	0	3755	120	0
2	A	34	0	0	1	0
2	B	37	0	0	1	0
2	C	34	0	0	2	0
2	D	35	0	0	0	0
All	All	15336	0	15020	433	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 14.

All (433) 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:365:LEU:HD12	1:B:395:ILE:HD11	1.31	1.11
1:B:365:LEU:HD13	1:B:386:PHE:HB3	1.49	0.93
1:A:120:VAL:HG12	1:A:124:MET:HE2	1.52	0.90
1:D:377:ARG:HG3	1:D:378:GLY:H	1.37	0.89
1:D:99:ARG:HG3	1:D:122:LEU:HD13	1.56	0.87
1:D:365:LEU:HD13	1:D:395:ILE:HD11	1.54	0.87
1:B:361:LYS:HE2	1:B:367:LEU:HD22	1.57	0.85
1:B:175:GLN:HG3	1:B:191:MET:HE1	1.59	0.83
1:D:272:LYS:HG3	1:D:307:ARG:HD2	1.59	0.82
1:B:365:LEU:HG	1:B:393:MET:SD	2.20	0.82
1:D:365:LEU:HD21	1:D:393:MET:SD	2.20	0.81
1:B:120:VAL:HG12	1:B:124:MET:HE2	1.63	0.80
1:A:365:LEU:HD21	1:A:393:MET:SD	2.21	0.80
1:D:365:LEU:HD11	1:D:393:MET:SD	2.22	0.80
1:D:159:VAL:H	1:D:187:ASN:HD21	1.31	0.79
1:C:365:LEU:HD22	1:C:389:LEU:HD22	1.64	0.79
1:A:159:VAL:HG12	1:A:187:ASN:ND2	1.98	0.79
1:C:175:GLN:HG3	1:C:191:MET:HE1	1.65	0.78
1:B:389:LEU:HD22	1:B:406:GLN:HB3	1.66	0.77
1:D:323:VAL:HG13	1:D:369:CYS:SG	2.26	0.75
1:B:389:LEU:HD12	1:B:393:MET:SD	2.26	0.75
1:D:359:SER:O	1:D:363:GLU:HG2	1.87	0.75
1:C:365:LEU:HD13	1:C:395:ILE:HD11	1.69	0.74
1:D:361:LYS:HE2	1:D:367:LEU:HD22	1.68	0.74
1:B:365:LEU:HD21	1:B:389:LEU:HA	1.69	0.74
1:A:386:PHE:HB3	1:A:389:LEU:HD21	1.70	0.73
1:A:359:SER:O	1:A:363:GLU:HG2	1.89	0.72
1:B:389:LEU:HD11	1:B:396:ALA:HB2	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:175:GLN:HG3	1:D:191:MET:HE1	1.71	0.71
1:A:365:LEU:HD11	1:A:393:MET:SD	2.29	0.71
1:D:251:HIS:O	1:D:255:VAL:HG23	1.92	0.70
1:D:377:ARG:HG3	1:D:378:GLY:N	2.07	0.69
1:D:120:VAL:HG12	1:D:124:MET:HE2	1.75	0.69
1:C:120:VAL:HG12	1:C:124:MET:HE2	1.75	0.69
1:C:77:ARG:HH11	1:C:77:ARG:HG2	1.56	0.68
1:C:365:LEU:HD21	1:C:393:MET:SD	2.33	0.68
1:A:365:LEU:CD2	1:A:393:MET:SD	2.81	0.68
1:B:292:PHE:HE2	1:B:457:ASP:HB2	1.56	0.68
1:A:241:VAL:HG23	1:A:263:LYS:HD3	1.77	0.67
1:C:159:VAL:H	1:C:187:ASN:HD21	1.41	0.67
1:C:365:LEU:HD21	1:C:389:LEU:HA	1.78	0.66
1:D:497:GLN:NE2	1:D:499:ASN:HD21	1.94	0.66
1:C:36:THR:OG1	1:C:50:HIS:HB3	1.95	0.66
1:D:159:VAL:N	1:D:187:ASN:HD21	1.92	0.66
1:B:102:LEU:HD21	1:B:203:TYR:HD2	1.61	0.65
1:B:413:MET:HE1	1:B:438:LYS:HG2	1.78	0.65
1:B:159:VAL:H	1:B:187:ASN:HD21	1.43	0.65
1:C:99:ARG:HD2	1:C:122:LEU:HB3	1.78	0.64
1:B:159:VAL:N	1:B:187:ASN:HD21	1.96	0.64
1:D:389:LEU:HD11	1:D:407:ILE:N	2.12	0.64
1:A:315:TYR:O	1:A:319:VAL:HG23	1.99	0.63
1:D:103:ALA:HB2	1:D:122:LEU:HD12	1.81	0.63
1:B:23:PHE:CZ	1:B:26:ASN:HA	2.33	0.62
1:D:365:LEU:HD13	1:D:389:LEU:HD23	1.81	0.62
1:C:395:ILE:HB	1:C:400:ILE:HD11	1.81	0.62
1:B:359:SER:O	1:B:363:GLU:HG2	2.00	0.62
1:C:495:VAL:HG22	1:C:496:PRO:HD2	1.80	0.62
1:B:171:PRO:HG3	1:B:197:THR:HG21	1.82	0.61
1:D:11:PRO:HB3	1:D:114:TYR:CZ	2.34	0.61
1:D:171:PRO:HG3	1:D:197:THR:HG21	1.82	0.61
1:A:365:LEU:HD11	1:A:389:LEU:HD13	1.83	0.61
1:D:389:LEU:HD13	1:D:406:GLN:HB3	1.83	0.61
1:B:102:LEU:HD21	1:B:203:TYR:CD2	2.35	0.60
1:B:76:TRP:CH2	1:B:84:ARG:HG2	2.36	0.60
1:C:361:LYS:HE2	1:C:367:LEU:HD22	1.83	0.60
1:A:365:LEU:CD1	1:A:393:MET:SD	2.89	0.60
1:C:294:LEU:CD1	1:C:405:MET:HA	2.31	0.60
1:A:238:VAL:O	1:A:261:ASN:ND2	2.35	0.60
1:C:161:VAL:HA	1:C:188:VAL:HG23	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:279:SER:HB3	1:D:311:GLN:HG2	1.84	0.59
1:A:365:LEU:HG	1:A:395:ILE:HD11	1.85	0.59
1:C:103:ALA:HB2	1:C:122:LEU:HD12	1.85	0.59
1:B:365:LEU:CD1	1:B:395:ILE:HD11	2.20	0.59
1:A:301:CYS:HG	1:A:303:CYS:HG	1.50	0.59
1:B:124:MET:HE3	1:B:173:LEU:HD22	1.84	0.59
1:D:365:LEU:HD21	1:D:389:LEU:HA	1.84	0.59
1:D:120:VAL:CG1	1:D:124:MET:HE2	2.32	0.58
1:A:349:GLN:O	1:A:353:VAL:HG23	2.03	0.58
1:A:120:VAL:CG1	1:A:124:MET:HE2	2.30	0.58
1:D:59:VAL:O	1:D:63:VAL:HG23	2.04	0.58
1:C:77:ARG:HG2	1:C:77:ARG:NH1	2.16	0.58
1:A:294:LEU:HD22	1:A:405:MET:HB2	1.85	0.57
1:A:490:THR:OG1	1:B:464:PRO:HG2	2.04	0.57
1:B:126:LEU:O	1:B:130:ARG:HB2	2.04	0.57
1:A:159:VAL:HG12	1:A:187:ASN:HD21	1.69	0.57
1:A:159:VAL:H	1:A:187:ASN:HD21	1.51	0.57
1:A:193:VAL:HG11	1:A:198:PRO:HA	1.86	0.57
1:A:262:LEU:HG	1:B:251:HIS:CE1	2.39	0.57
1:A:371:GLY:HA2	1:A:384:THR:OG1	2.05	0.57
1:D:197:THR:O	1:D:197:THR:HG23	2.04	0.57
1:C:359:SER:O	1:C:363:GLU:HG2	2.05	0.56
1:A:272:LYS:HG3	1:A:307:ARG:HD2	1.86	0.56
1:C:99:ARG:HG3	1:C:122:LEU:HD13	1.86	0.56
1:D:70:PHE:CZ	1:D:158:PRO:HB2	2.41	0.56
1:D:300:GLN:HE22	1:D:345:VAL:H	1.52	0.56
1:D:389:LEU:HD11	1:D:406:GLN:CA	2.35	0.56
1:C:23:PHE:CZ	1:C:26:ASN:HA	2.40	0.56
1:A:365:LEU:HD11	1:A:389:LEU:HA	1.87	0.56
1:A:365:LEU:HD12	1:A:389:LEU:HD22	1.87	0.56
1:B:309:PHE:CE1	1:B:408:LEU:HD22	2.41	0.56
1:C:15:PRO:HD2	1:C:108:LEU:HD22	1.87	0.56
1:B:55:ASP:CG	1:B:227:THR:HG23	2.31	0.56
1:C:443:SER:HA	1:C:451:VAL:HG11	1.87	0.56
1:A:361:LYS:HD3	1:A:367:LEU:HD22	1.88	0.55
1:C:26:ASN:O	1:C:209:LYS:HE2	2.05	0.55
1:A:323:VAL:HG13	1:A:369:CYS:SG	2.46	0.55
1:C:171:PRO:HG3	1:C:197:THR:HG21	1.88	0.55
1:C:20:ASN:HA	1:C:202:LEU:HD13	1.89	0.55
1:B:365:LEU:HD11	1:B:389:LEU:HD13	1.88	0.55
1:D:365:LEU:CD2	1:D:393:MET:SD	2.95	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:365:LEU:CD2	1:D:389:LEU:HA	2.37	0.54
1:A:126:LEU:O	1:A:130:ARG:HB2	2.07	0.54
1:B:170:PHE:HE1	2:B:513:HOH:O	1.89	0.54
1:D:389:LEU:O	1:D:408:LEU:HD23	2.07	0.54
1:A:107:THR:HG23	1:A:112:LYS:O	2.07	0.54
1:C:164:GLN:CD	1:C:178:LYS:HB3	2.33	0.54
1:C:245:GLY:O	1:C:269:ILE:HG22	2.08	0.54
1:A:15:PRO:HD2	1:A:108:LEU:HD22	1.89	0.53
1:B:365:LEU:HD22	1:B:387:GLY:O	2.08	0.53
1:B:192:LYS:HE2	1:B:223:GLY:C	2.34	0.53
1:D:160:GLY:H	1:D:187:ASN:HD22	1.56	0.53
1:B:11:PRO:HB3	1:B:114:TYR:CE1	2.44	0.53
1:D:70:PHE:HZ	1:D:158:PRO:HB2	1.74	0.53
1:A:198:PRO:O	1:A:202:LEU:HG	2.08	0.53
1:C:349:GLN:O	1:C:353:VAL:HG23	2.09	0.53
1:C:460:GLY:HA3	1:D:146:ILE:HG13	1.91	0.53
1:C:371:GLY:HA2	1:C:384:THR:OG1	2.08	0.53
1:A:197:THR:HG23	1:A:197:THR:O	2.07	0.53
1:B:59:VAL:O	1:B:63:VAL:HG23	2.09	0.53
1:D:40:VAL:HG12	1:D:41:ASN:N	2.24	0.52
1:A:443:SER:HA	1:A:451:VAL:HG11	1.91	0.52
1:D:347:GLU:O	1:D:350:PHE:HB3	2.10	0.52
1:D:365:LEU:CD1	1:D:393:MET:SD	2.96	0.52
1:D:329:ARG:HE	1:D:341:GLN:HE21	1.58	0.52
1:A:148:GLY:O	1:A:498:LYS:HD3	2.09	0.52
1:B:268:GLU:OE1	1:B:476:GLU:HG3	2.09	0.52
1:C:294:LEU:HD11	1:C:405:MET:HA	1.92	0.52
1:B:391:ASP:OD2	1:B:419:ARG:HD2	2.09	0.52
1:D:11:PRO:HB3	1:D:114:TYR:CE1	2.44	0.52
1:A:400:ILE:HB	2:A:524:HOH:O	2.09	0.52
1:D:277:ILE:HD12	1:D:277:ILE:N	2.26	0.52
1:C:424:LYS:HB3	1:C:470:LEU:HD12	1.92	0.51
1:A:99:ARG:HG2	1:A:118:TYR:CE2	2.46	0.51
1:A:311:GLN:NE2	1:A:411:LYS:O	2.43	0.51
1:B:235:HIS:HB3	1:B:238:VAL:HG23	1.93	0.51
1:B:120:VAL:CG1	1:B:124:MET:HE2	2.39	0.51
1:B:371:GLY:HA2	1:B:384:THR:OG1	2.11	0.51
1:C:99:ARG:NH1	1:C:118:TYR:O	2.44	0.51
1:A:11:PRO:HB3	1:A:114:TYR:CE1	2.46	0.51
1:C:358:LYS:O	1:C:362:GLU:HG3	2.11	0.51
1:D:23:PHE:CZ	1:D:26:ASN:HA	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:THR:HG23	1:C:48:ILE:HB	1.93	0.51
1:A:103:ALA:HB2	1:A:122:LEU:HD12	1.94	0.50
1:D:121:ASP:O	1:D:125:VAL:HG23	2.11	0.50
1:A:11:PRO:HB3	1:A:114:TYR:CZ	2.46	0.50
1:B:227:THR:HG22	1:B:228:ALA:N	2.25	0.50
1:B:241:VAL:CG1	1:B:265:VAL:HG22	2.41	0.50
1:B:77:ARG:HG2	1:B:77:ARG:HH11	1.77	0.50
1:B:409:LYS:O	1:B:419:ARG:NH2	2.43	0.50
1:C:120:VAL:O	1:C:124:MET:HG3	2.12	0.50
1:D:224:PHE:HB2	1:D:227:THR:OG1	2.12	0.50
1:A:193:VAL:CG1	1:A:198:PRO:HA	2.41	0.50
1:D:365:LEU:CD1	1:D:395:ILE:HD11	2.34	0.50
1:A:227:THR:HG22	1:A:228:ALA:N	2.25	0.50
1:B:269:ILE:HG13	1:B:471:SER:HA	1.94	0.50
1:B:198:PRO:O	1:B:202:LEU:HG	2.11	0.50
1:C:70:PHE:O	1:C:77:ARG:HD2	2.12	0.50
1:C:413:MET:HA	1:C:416:VAL:HG12	1.94	0.50
1:C:497:GLN:NE2	1:C:499:ASN:HD21	2.08	0.50
1:D:319:VAL:O	1:D:323:VAL:HG23	2.11	0.50
1:A:121:ASP:O	1:A:125:VAL:HG23	2.12	0.50
1:A:309:PHE:HB3	1:A:410:PHE:CD2	2.47	0.50
1:C:291:HIS:NE2	1:C:329:ARG:NH1	2.59	0.50
1:B:183:LEU:HD13	1:B:213:PHE:CE2	2.47	0.49
1:C:489:LYS:HB2	1:D:468:TYR:OH	2.12	0.49
1:D:497:GLN:HE21	1:D:499:ASN:HD21	1.60	0.49
1:B:23:PHE:CE1	1:B:26:ASN:HA	2.48	0.49
1:D:21:GLN:HB3	1:D:29:HIS:O	2.12	0.49
1:D:245:GLY:O	1:D:269:ILE:HG22	2.11	0.49
1:D:292:PHE:CE2	1:D:457:ASP:HB2	2.48	0.49
1:B:11:PRO:HB3	1:B:114:TYR:CZ	2.47	0.49
1:B:208:ILE:HD13	1:B:218:VAL:HG11	1.93	0.49
1:B:413:MET:O	1:B:417:VAL:HG23	2.11	0.49
1:C:175:GLN:CG	1:C:191:MET:HE1	2.39	0.49
1:A:168:TRP:HA	1:A:197:THR:HG21	1.95	0.49
1:D:365:LEU:CG	1:D:393:MET:SD	3.00	0.49
1:A:278:MET:HE2	1:A:438:LYS:HE3	1.93	0.49
1:B:294:LEU:CD1	1:B:405:MET:HA	2.42	0.49
1:B:349:GLN:O	1:B:353:VAL:HG23	2.12	0.49
1:C:198:PRO:O	1:C:202:LEU:HG	2.13	0.49
1:C:336:ASP:OD2	1:C:338:ARG:HD2	2.12	0.49
1:C:251:HIS:HA	1:D:262:LEU:HD11	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:389:LEU:CD2	1:B:406:GLN:HB3	2.41	0.49
1:C:11:PRO:HB3	1:C:114:TYR:CZ	2.48	0.49
1:A:300:GLN:HE22	1:A:345:VAL:H	1.59	0.48
1:C:227:THR:HG22	1:C:228:ALA:N	2.27	0.48
1:C:279:SER:HA	1:C:314:ILE:HD13	1.94	0.48
1:B:168:TRP:HA	1:B:197:THR:HG21	1.95	0.48
1:C:413:MET:HE3	1:C:441:TYR:CD2	2.48	0.48
1:A:270:GLY:O	1:A:471:SER:HB2	2.13	0.48
1:C:255:VAL:HG13	1:D:255:VAL:HG13	1.95	0.48
1:A:70:PHE:CZ	1:A:158:PRO:HB2	2.47	0.48
1:A:171:PRO:HG3	1:A:197:THR:HG21	1.95	0.48
1:A:294:LEU:HD22	1:A:405:MET:CB	2.43	0.48
1:D:389:LEU:HD11	1:D:407:ILE:H	1.77	0.48
1:C:294:LEU:HD13	1:C:405:MET:HA	1.96	0.48
1:A:390:GLN:O	1:A:393:MET:HG3	2.13	0.48
1:A:261:ASN:O	1:A:262:LEU:HB2	2.14	0.48
1:C:183:LEU:HD13	1:C:213:PHE:CE2	2.49	0.48
1:D:99:ARG:HG2	1:D:118:TYR:CE2	2.49	0.48
1:A:40:VAL:HG12	1:A:41:ASN:N	2.28	0.48
1:C:42:PRO:HB3	1:C:345:VAL:O	2.14	0.48
1:C:172:LEU:HD21	1:C:200:THR:HB	1.96	0.48
1:D:126:LEU:O	1:D:130:ARG:HB2	2.14	0.48
1:D:272:LYS:CG	1:D:307:ARG:HD2	2.40	0.48
1:A:185:THR:HG21	1:A:485:TYR:O	2.14	0.47
1:C:389:LEU:HD12	1:C:396:ALA:HB2	1.96	0.47
1:D:430:ALA:HB2	1:D:456:TYR:CD1	2.49	0.47
1:C:319:VAL:O	1:C:323:VAL:HG23	2.13	0.47
1:D:62:ALA:CB	1:D:221:ILE:HD11	2.44	0.47
1:D:159:VAL:HG11	1:D:240:LYS:HB2	1.95	0.47
1:A:188:VAL:HG13	1:A:217:VAL:HA	1.95	0.47
1:B:195:GLU:HB3	1:B:223:GLY:O	2.13	0.47
1:C:365:LEU:HD11	1:C:393:MET:SD	2.54	0.47
1:C:317:GLU:O	1:C:321:ARG:HG2	2.14	0.47
1:C:365:LEU:CG	1:C:393:MET:SD	3.02	0.47
1:B:112:LYS:HE2	1:B:297:ASN:OD1	2.14	0.47
1:A:446:LEU:O	1:B:489:LYS:NZ	2.41	0.47
1:A:467:GLY:O	1:A:475:ARG:NH2	2.48	0.47
1:C:11:PRO:HB3	1:C:114:TYR:CE1	2.50	0.47
1:C:170:PHE:HB3	1:C:173:LEU:HB3	1.97	0.47
1:D:389:LEU:CD1	1:D:407:ILE:N	2.76	0.47
1:A:195:GLU:HB3	1:A:223:GLY:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:316:ALA:O	1:D:320:GLU:HG2	2.15	0.47
1:D:404:VAL:HG12	1:D:406:GLN:OE1	2.15	0.47
1:B:70:PHE:CZ	1:B:158:PRO:HB2	2.50	0.47
1:A:247:THR:HA	1:A:269:ILE:HD12	1.96	0.46
1:D:365:LEU:HD22	1:D:389:LEU:HD23	1.96	0.46
1:A:462:GLN:O	1:B:144:ILE:HG21	2.14	0.46
1:B:424:LYS:O	1:B:469:LYS:HB2	2.15	0.46
1:D:235:HIS:HB3	1:D:238:VAL:HG23	1.96	0.46
1:A:386:PHE:CB	1:A:389:LEU:HD21	2.40	0.46
1:B:365:LEU:HD23	1:B:365:LEU:HA	1.55	0.46
1:A:451:VAL:HG23	1:B:489:LYS:HD2	1.97	0.46
1:D:389:LEU:CD1	1:D:406:GLN:HB3	2.46	0.46
1:A:251:HIS:O	1:A:255:VAL:HG23	2.16	0.46
1:B:467:GLY:O	1:B:475:ARG:NH2	2.49	0.46
1:C:157:GLU:OE2	1:C:489:LYS:HE2	2.16	0.46
1:C:365:LEU:CD2	1:C:393:MET:SD	3.01	0.46
1:D:389:LEU:HD11	1:D:406:GLN:HA	1.95	0.46
1:A:279:SER:HB3	1:A:311:GLN:HG2	1.98	0.46
1:D:268:GLU:OE2	1:D:466:GLY:HA2	2.15	0.46
1:C:205:ALA:HA	1:C:208:ILE:HD12	1.98	0.46
1:D:266:THR:O	1:D:267:LEU:HD23	2.16	0.46
1:A:429:ALA:HB1	1:A:446:LEU:HD13	1.97	0.46
1:D:424:LYS:O	1:D:469:LYS:HB2	2.15	0.46
1:C:124:MET:HE3	1:C:173:LEU:HD22	1.97	0.45
1:C:389:LEU:HB3	1:C:408:LEU:HG	1.97	0.45
1:D:12:ASN:O	1:D:15:PRO:HD3	2.15	0.45
1:D:413:MET:HA	1:D:416:VAL:HG12	1.98	0.45
1:A:283:MET:HE1	1:A:318:PHE:HA	1.98	0.45
1:A:488:VAL:O	1:B:475:ARG:NH1	2.49	0.45
1:B:395:ILE:HB	1:B:400:ILE:HD11	1.98	0.45
1:C:135:TRP:CG	1:C:482:LEU:HD11	2.50	0.45
1:D:294:LEU:HD12	1:D:306:SER:HA	1.98	0.45
1:D:497:GLN:HE21	1:D:499:ASN:ND2	2.14	0.45
1:A:23:PHE:CZ	1:A:26:ASN:HA	2.52	0.45
1:A:55:ASP:CG	1:A:227:THR:HG23	2.41	0.45
1:A:272:LYS:HG3	1:A:307:ARG:CD	2.47	0.45
1:D:306:SER:O	1:D:406:GLN:HB2	2.17	0.45
1:A:42:PRO:HG3	1:A:110:ASN:O	2.17	0.45
1:B:317:GLU:O	1:B:321:ARG:HG3	2.17	0.45
1:C:44:THR:HA	1:C:377:ARG:HD2	1.99	0.45
1:D:70:PHE:CE2	1:D:160:GLY:HA2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:24:ILE:HD12	1:C:62:ALA:HB2	1.99	0.45
1:C:293:ALA:HA	1:C:456:TYR:CD2	2.52	0.45
1:B:389:LEU:HD22	1:B:406:GLN:CB	2.42	0.44
1:C:84:ARG:NH1	1:C:184:ALA:O	2.50	0.44
1:D:272:LYS:HG3	1:D:307:ARG:CD	2.38	0.44
1:D:283:MET:O	1:D:287:VAL:HG23	2.16	0.44
1:A:283:MET:CE	1:A:321:ARG:HD2	2.47	0.44
1:C:159:VAL:H	1:C:187:ASN:ND2	2.12	0.44
1:C:269:ILE:HB	1:C:270:GLY:H	1.69	0.44
1:D:107:THR:HG23	1:D:112:LYS:O	2.17	0.44
1:A:185:THR:HG22	1:A:185:THR:O	2.17	0.44
1:C:126:LEU:O	1:C:130:ARG:HB2	2.18	0.44
1:C:389:LEU:HG	1:C:406:GLN:HB3	2.00	0.44
1:D:33:SER:OG	1:D:35:LYS:HG3	2.18	0.44
1:D:371:GLY:HA2	1:D:384:THR:OG1	2.17	0.44
1:C:241:VAL:HG23	1:C:263:LYS:HG3	2.00	0.44
1:D:241:VAL:CG2	1:D:263:LYS:HE2	2.47	0.44
1:A:113:PRO:HB2	1:A:116:ILE:HG12	2.00	0.44
1:B:294:LEU:HD12	1:B:306:SER:HA	2.00	0.44
1:C:365:LEU:HD13	1:C:395:ILE:CD1	2.45	0.44
1:C:365:LEU:HG	1:C:393:MET:SD	2.56	0.44
1:D:80:ASP:HB2	1:D:83:GLU:HG2	1.99	0.44
1:B:39:THR:HG23	1:B:48:ILE:HB	1.99	0.44
1:B:389:LEU:HD23	1:B:408:LEU:HD11	1.99	0.44
1:C:404:VAL:HG12	1:C:406:GLN:OE1	2.18	0.44
1:C:424:LYS:O	1:C:469:LYS:HB2	2.17	0.44
1:D:170:PHE:HB3	1:D:173:LEU:HB3	1.99	0.44
1:B:306:SER:O	1:B:406:GLN:HB2	2.18	0.44
1:B:389:LEU:HD21	1:B:396:ALA:HB2	1.99	0.44
1:C:205:ALA:O	1:C:208:ILE:HB	2.18	0.44
1:D:295:PHE:HE2	1:D:405:MET:HE3	1.83	0.44
1:B:160:GLY:H	1:B:187:ASN:ND2	2.16	0.44
1:B:463:SER:O	1:B:477:LEU:HB2	2.18	0.44
1:C:410:PHE:CD2	1:C:416:VAL:HB	2.53	0.44
1:D:161:VAL:HA	1:D:188:VAL:HG23	2.00	0.44
1:D:389:LEU:HD11	1:D:406:GLN:C	2.43	0.44
1:B:292:PHE:CE2	1:B:457:ASP:HB2	2.45	0.43
1:D:161:VAL:HA	1:D:188:VAL:CG2	2.48	0.43
1:B:42:PRO:HB2	1:B:343:PRO:HG2	2.00	0.43
1:B:251:HIS:O	1:B:255:VAL:HG23	2.18	0.43
1:B:294:LEU:HD13	1:B:405:MET:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298:GLN:HG2	1:B:343:PRO:O	2.18	0.43
1:B:315:TYR:CD2	1:B:409:LYS:HE2	2.53	0.43
1:C:21:GLN:HB3	1:C:29:HIS:O	2.17	0.43
1:C:55:ASP:CG	1:C:227:THR:HG23	2.43	0.43
1:C:498:LYS:C	1:C:498:LYS:HD2	2.43	0.43
1:D:244:THR:HA	1:D:268:GLU:O	2.19	0.43
1:B:42:PRO:HG3	1:B:110:ASN:O	2.19	0.43
1:C:77:ARG:HH11	1:C:77:ARG:CG	2.29	0.43
1:A:365:LEU:HD13	1:A:365:LEU:HA	1.75	0.43
1:A:144:ILE:HG21	1:B:462:GLN:O	2.18	0.43
1:C:365:LEU:HG	1:C:393:MET:CE	2.49	0.43
1:D:149:ASP:C	1:D:495:VAL:HG12	2.44	0.43
1:D:443:SER:HA	1:D:451:VAL:HG11	2.01	0.43
1:A:155:ARG:NH2	1:B:444:GLN:HG3	2.34	0.43
1:A:413:MET:O	1:A:417:VAL:HG23	2.19	0.43
1:A:500:SER:OXT	1:D:158:PRO:HD3	2.18	0.43
1:B:121:ASP:O	1:B:125:VAL:HG23	2.19	0.43
1:B:333:ASN:HA	1:B:334:PRO:HD3	1.90	0.43
1:C:389:LEU:CB	1:C:408:LEU:HG	2.49	0.43
1:D:249:VAL:HA	1:D:252:LEU:HD12	2.00	0.43
1:B:168:TRP:HZ2	1:B:349:GLN:NE2	2.16	0.43
1:D:40:VAL:CG1	1:D:41:ASN:N	2.81	0.43
1:C:39:THR:HG21	1:C:199:LEU:HD11	2.01	0.43
1:D:386:PHE:HB3	1:D:389:LEU:HD21	2.01	0.43
1:A:99:ARG:HD3	1:A:122:LEU:HB3	2.01	0.42
1:A:448:ALA:O	1:B:489:LYS:HD3	2.19	0.42
1:B:279:SER:CB	1:B:311:GLN:HG2	2.48	0.42
1:D:410:PHE:CD2	1:D:416:VAL:HB	2.54	0.42
1:A:40:VAL:CG1	1:A:41:ASN:N	2.82	0.42
1:A:424:LYS:HD2	1:A:470:LEU:CD1	2.49	0.42
1:C:99:ARG:HG2	1:C:118:TYR:CE1	2.54	0.42
1:B:12:ASN:O	1:B:15:PRO:HD3	2.18	0.42
1:A:244:THR:HA	1:A:268:GLU:O	2.20	0.42
1:C:92:ALA:CB	1:C:130:ARG:HD3	2.50	0.42
1:C:240:LYS:HG2	1:C:241:VAL:N	2.34	0.42
1:C:116:ILE:HG21	2:C:512:HOH:O	2.19	0.42
1:C:272:LYS:HD2	1:C:272:LYS:HA	1.88	0.42
1:D:146:ILE:HG12	1:D:147:ASP:N	2.33	0.42
1:A:347:GLU:O	1:A:350:PHE:HB3	2.19	0.42
1:A:413:MET:HA	1:A:416:VAL:HG12	2.02	0.42
1:B:208:ILE:CD1	1:B:218:VAL:HG11	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ILE:HD12	1:B:62:ALA:HB2	2.01	0.42
1:B:175:GLN:CG	1:B:191:MET:HE1	2.41	0.42
1:B:300:GLN:HE22	1:B:345:VAL:H	1.68	0.42
1:B:365:LEU:HD11	1:B:389:LEU:CD1	2.48	0.42
1:B:365:LEU:HB3	1:B:386:PHE:CD1	2.54	0.42
1:A:21:GLN:HB3	1:A:29:HIS:O	2.20	0.42
1:B:117:SER:HA	1:B:121:ASP:HB2	2.01	0.42
1:D:185:THR:O	1:D:185:THR:HG22	2.19	0.42
1:D:202:LEU:O	1:D:205:ALA:HB3	2.20	0.42
1:B:244:THR:HG23	1:B:268:GLU:HB3	2.02	0.42
1:B:244:THR:HA	1:B:268:GLU:O	2.20	0.42
1:B:279:SER:HB3	1:B:311:GLN:HG2	2.01	0.42
1:C:106:GLU:O	1:C:110:ASN:HB3	2.20	0.42
1:C:115:ILE:HG12	2:C:524:HOH:O	2.19	0.42
1:D:395:ILE:HB	1:D:400:ILE:HD11	2.02	0.42
1:A:59:VAL:O	1:A:63:VAL:HG23	2.20	0.42
1:C:303:CYS:SG	1:C:459:PHE:HZ	2.43	0.42
1:D:309:PHE:CZ	1:D:408:LEU:HD13	2.55	0.42
1:A:154:THR:HA	1:A:489:LYS:O	2.20	0.41
1:C:44:THR:HA	1:C:377:ARG:CD	2.50	0.41
1:C:193:VAL:HG11	1:C:198:PRO:HA	2.02	0.41
1:D:106:GLU:O	1:D:110:ASN:HB3	2.20	0.41
1:A:55:ASP:OD1	1:A:227:THR:HG23	2.20	0.41
1:A:170:PHE:HB3	1:A:173:LEU:HB3	2.02	0.41
1:C:146:ILE:HA	1:D:462:GLN:CG	2.50	0.41
1:C:168:TRP:CB	1:C:197:THR:HG22	2.49	0.41
1:C:310:VAL:HG21	1:C:318:PHE:CD2	2.56	0.41
1:D:331:VAL:HG22	1:D:341:GLN:HB3	2.02	0.41
1:D:462:GLN:H	1:D:462:GLN:CD	2.27	0.41
1:B:135:TRP:CD1	1:D:138:LYS:HE3	2.55	0.41
1:B:495:VAL:HB	1:B:496:PRO:HD2	2.03	0.41
1:C:430:ALA:HB2	1:C:456:TYR:CD1	2.55	0.41
1:D:413:MET:O	1:D:417:VAL:HG23	2.21	0.41
1:A:389:LEU:HD11	1:A:395:ILE:HD11	2.02	0.41
1:A:430:ALA:HA	1:A:452:TRP:O	2.21	0.41
1:C:214:PRO:HA	1:C:215:PRO:HD3	1.95	0.41
1:C:269:ILE:HG13	1:C:471:SER:C	2.46	0.41
1:A:135:TRP:CG	1:A:482:LEU:HD11	2.56	0.41
1:A:270:GLY:O	1:A:471:SER:CB	2.69	0.41
1:A:464:PRO:HG2	1:B:490:THR:OG1	2.21	0.41
1:B:99:ARG:HG2	1:B:118:TYR:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:LYS:HE2	1:C:297:ASN:OD1	2.19	0.41
1:D:24:ILE:HD12	1:D:61:ARG:CZ	2.50	0.41
1:A:247:THR:O	1:A:251:HIS:HD2	2.04	0.41
1:B:40:VAL:HG12	1:B:41:ASN:N	2.36	0.41
1:B:185:THR:O	1:B:185:THR:HG22	2.21	0.41
1:A:12:ASN:O	1:A:15:PRO:HD3	2.21	0.41
1:B:241:VAL:HG12	1:B:265:VAL:HG22	2.02	0.41
1:B:377:ARG:HH11	1:B:377:ARG:HG3	1.86	0.41
1:B:461:ALA:HA	1:B:477:LEU:HD22	2.01	0.41
1:D:33:SER:O	1:D:34:LYS:HB2	2.19	0.41
1:C:195:GLU:HB3	1:C:223:GLY:O	2.21	0.41
1:D:33:SER:OG	1:D:35:LYS:HE3	2.21	0.41
1:B:77:ARG:HG2	1:B:77:ARG:NH1	2.36	0.40
1:B:154:THR:HA	1:B:489:LYS:O	2.21	0.40
1:B:391:ASP:CG	1:B:419:ARG:HH11	2.28	0.40
1:D:278:MET:HE2	1:D:413:MET:SD	2.61	0.40
1:D:430:ALA:HA	1:D:452:TRP:O	2.21	0.40
1:A:208:ILE:HD13	1:A:218:VAL:HG11	2.04	0.40
1:A:365:LEU:CD1	1:A:389:LEU:HD22	2.51	0.40
1:D:157:GLU:OE2	1:D:489:LYS:HE2	2.22	0.40
1:C:13:GLN:HE22	1:C:335:PHE:HB3	1.86	0.40
1:C:41:ASN:ND2	1:C:108:LEU:O	2.55	0.40
1:C:468:TYR:OH	1:D:489:LYS:HB2	2.22	0.40
1:D:463:SER:O	1:D:477:LEU:HB2	2.21	0.40
1:A:63:VAL:HG11	1:A:235:HIS:CE1	2.57	0.40
1:B:197:THR:O	1:B:197:THR:HG23	2.20	0.40
1:B:262:LEU:HD12	1:B:262:LEU:HA	1.94	0.40
1:C:244:THR:HA	1:C:268:GLU:O	2.21	0.40
1:C:365:LEU:CD2	1:C:389:LEU:HA	2.49	0.40
1:D:365:LEU:HA	1:D:365:LEU:HD23	1.78	0.40
1:D:469:LYS:C	1:D:471:SER:H	2.30	0.40
1:A:255:VAL:HG13	1:B:255:VAL:HG13	2.02	0.40
1:A:372:GLY:O	1:A:382:GLN:HG3	2.22	0.40
1:B:141:GLY:HA3	1:C:143:THR:OG1	2.22	0.40
1:B:146:ILE:HG12	1:B:147:ASP:N	2.36	0.40
1:D:160:GLY:H	1:D:187:ASN:ND2	2.19	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	491/499 (98%)	464 (94%)	26 (5%)	1 (0%)	43	60
1	B	491/499 (98%)	464 (94%)	26 (5%)	1 (0%)	43	60
1	C	491/499 (98%)	463 (94%)	27 (6%)	1 (0%)	43	60
1	D	491/499 (98%)	466 (95%)	24 (5%)	1 (0%)	43	60
All	All	1964/1996 (98%)	1857 (95%)	103 (5%)	4 (0%)	43	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	ILE
1	B	24	ILE
1	C	24	ILE
1	D	24	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	398/401 (99%)	381 (96%)	17 (4%)	26	44
1	B	398/401 (99%)	382 (96%)	16 (4%)	28	47
1	C	398/401 (99%)	379 (95%)	19 (5%)	23	39
1	D	398/401 (99%)	386 (97%)	12 (3%)	36	57
All	All	1592/1604 (99%)	1528 (96%)	64 (4%)	28	47

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	29	HIS
1	A	64	LYS
1	A	99	ARG
1	A	130	ARG
1	A	264	ARG
1	A	269	ILE
1	A	294	LEU
1	A	311	GLN
1	A	338	ARG
1	A	365	LEU
1	A	377	ARG
1	A	401	PHE
1	A	422	ASN
1	A	468	TYR
1	A	470	LEU
1	A	475	ARG
1	B	99	ARG
1	B	122	LEU
1	B	130	ARG
1	B	262	LEU
1	B	263	LYS
1	B	264	ARG
1	B	268	GLU
1	B	269	ILE
1	B	279	SER
1	B	306	SER
1	B	311	GLN
1	B	327	LYS
1	B	365	LEU
1	B	377	ARG
1	B	468	TYR
1	B	475	ARG
1	C	40	VAL
1	C	99	ARG
1	C	130	ARG
1	C	159	VAL
1	C	241	VAL
1	C	262	LEU
1	C	263	LYS
1	C	264	ARG
1	C	266	THR

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Mol	Chain	Res	Type
1	C	269	ILE
1	C	337	SER
1	C	338	ARG
1	C	348	THR
1	C	369	CYS
1	C	377	ARG
1	C	394	THR
1	C	401	PHE
1	C	475	ARG
1	C	495	VAL
1	D	64	LYS
1	D	90	ARG
1	D	130	ARG
1	D	262	LEU
1	D	269	ILE
1	D	288	GLU
1	D	311	GLN
1	D	321	ARG
1	D	389	LEU
1	D	401	PHE
1	D	475	ARG
1	D	476	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	14	GLN
1	A	26	ASN
1	A	71	GLN
1	A	140	HIS
1	A	175	GLN
1	A	187	ASN
1	A	261	ASN
1	A	275	ASN
1	A	462	GLN
1	B	13	GLN
1	B	25	ASN
1	B	26	ASN
1	B	41	ASN
1	B	50	HIS
1	B	187	ASN

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Mol	Chain	Res	Type
1	B	300	GLN
1	B	311	GLN
1	B	422	ASN
1	B	447	GLN
1	C	26	ASN
1	C	41	ASN
1	C	89	ASN
1	C	187	ASN
1	C	275	ASN
1	C	382	GLN
1	C	422	ASN
1	C	462	GLN
1	C	497	GLN
1	D	26	ASN
1	D	187	ASN
1	D	275	ASN
1	D	300	GLN
1	D	341	GLN
1	D	344	GLN
1	D	422	ASN
1	D	447	GLN
1	D	497	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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	493/499 (98%)	-0.52	1 (0%) 91 90	2, 10, 24, 42	0
1	B	493/499 (98%)	-0.49	3 (0%) 85 83	2, 9, 24, 35	0
1	C	493/499 (98%)	-0.28	2 (0%) 88 87	3, 21, 40, 53	0
1	D	493/499 (98%)	-0.36	2 (0%) 88 87	2, 14, 32, 53	0
All	All	1972/1996 (98%)	-0.41	8 (0%) 88 87	2, 12, 32, 53	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	370	GLY	3.3
1	C	74	SER	3.3
1	A	74	SER	2.6
1	B	197	THR	2.6
1	C	338	ARG	2.4
1	B	261	ASN	2.4
1	D	376	ASP	2.1
1	B	363	GLU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.