



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:53 AM UTC

PDB ID : 1K5H / pdb_00001k5h
Title : 1-deoxy-D-xylulose-5-phosphate reductoisomerase
Authors : Reuter, K.; Sanderbrand, S.; Jomaa, H.; Wiesner, J.; Steinbrecher, I.; Beck, E.; Hintz, M.; Klebe, G.; Stubbs, M.T.
Deposited on : 2001-10-10
Resolution : 2.50 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ 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

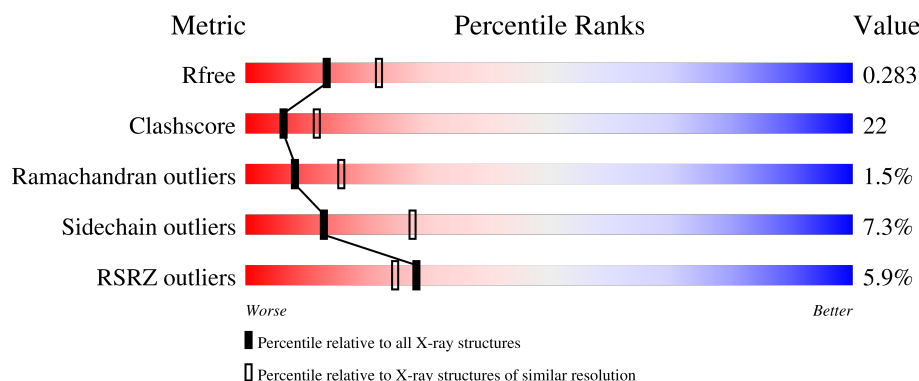
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	398	4% 59% 35% 6%
1	B	398	11% 56% 36% 6%
1	C	398	3% 58% 38%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9126 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 1-deoxy-D-xylulose-5-phosphate reductoisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	398	Total	C	N	O	S	0	0	0
			3030	1892	534	577	27			
1	B	391	Total	C	N	O	S	0	0	0
			2978	1860	526	567	25			
1	C	398	Total	C	N	O	S	0	0	0
			3030	1892	534	577	27			

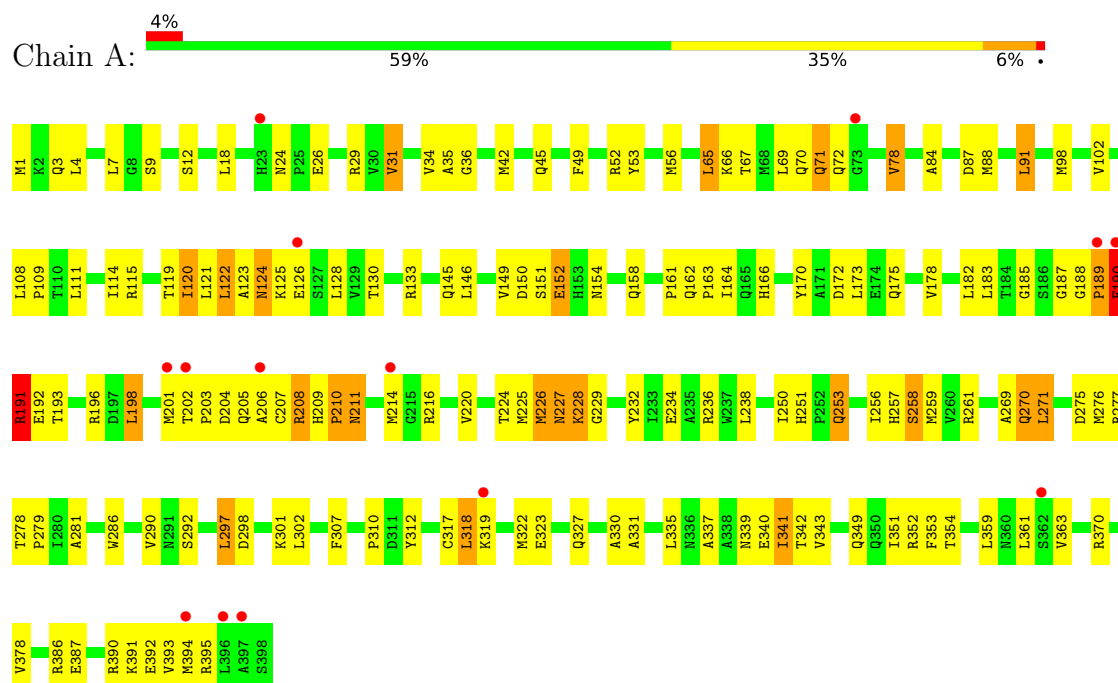
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	23	Total	O	0	0
			23	23		
2	B	25	Total	O	0	0
			25	25		
2	C	40	Total	O	0	0
			40	40		

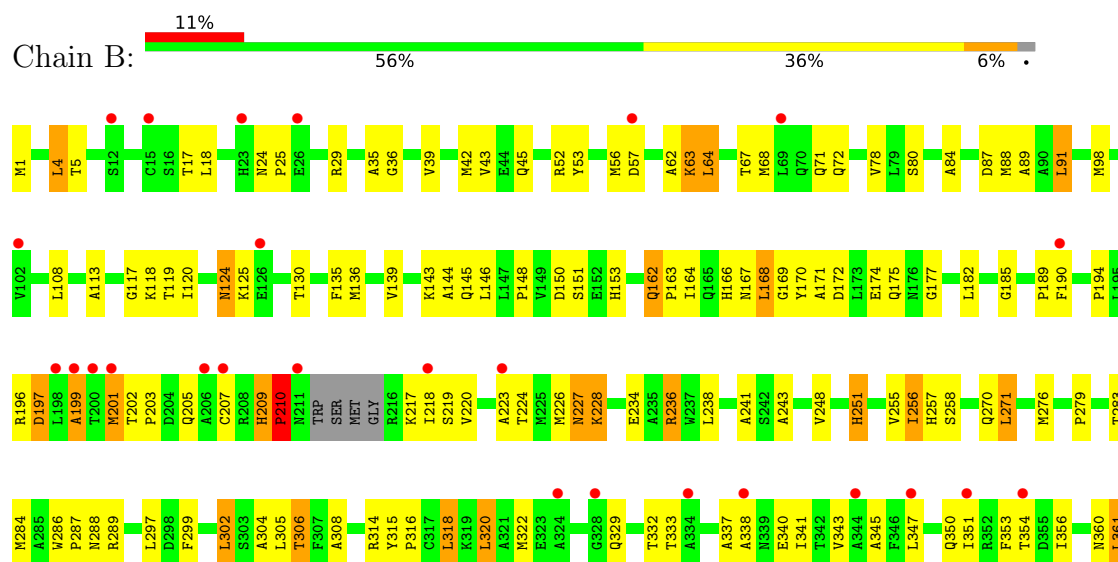
3 Residue-property plots

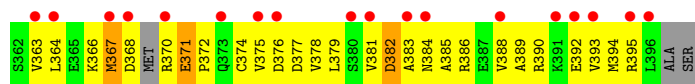
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 1-deoxy-D-xylulose-5-phosphate reductoisomerase

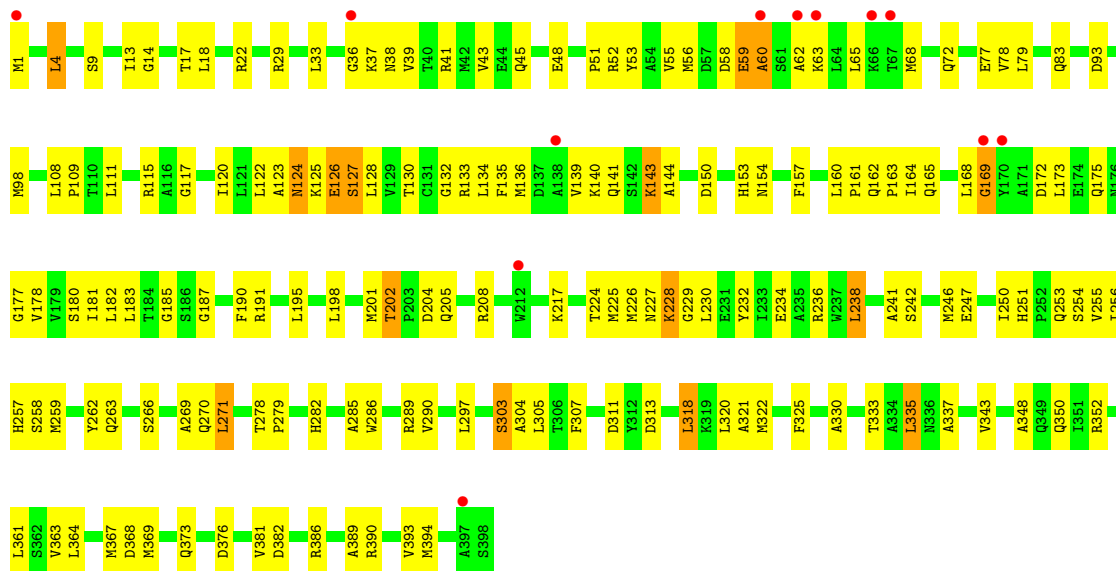


- Molecule 1: 1-deoxy-D-xylulose-5-phosphate reductoisomerase





● Molecule 1: 1-deoxy-D-xylulose-5-phosphate reductoisomerase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	101.27Å 249.26Å 132.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.45 – 2.50 33.45 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.6 (33.45-2.50) 90.6 (33.45-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.232 , 0.284 0.233 , 0.283	Depositor DCC
R_{free} test set	5816 reflections (10.16%)	wwPDB-VP
Wilson B-factor (Å ²)	49.4	Xtriage
Anisotropy	0.576	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.8	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	0.94	EDS
Total number of atoms	9126	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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	0.46	1/3078 (0.0%)	0.95	12/4171 (0.3%)
1	B	0.46	1/3022 (0.0%)	1.00	11/4094 (0.3%)
1	C	0.50	0/3078	0.98	8/4171 (0.2%)
All	All	0.47	2/9178 (0.0%)	0.98	31/12436 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	192	GLU	N-CA	5.46	1.52	1.45
1	B	209	HIS	CB-CG	5.23	1.57	1.50

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	210	PRO	CA-N-CD	-15.18	90.75	112.00
1	C	255	VAL	N-CA-C	-7.82	103.24	110.82
1	A	251	HIS	CA-C-N	7.79	129.57	119.84
1	A	251	HIS	C-N-CA	7.79	129.57	119.84
1	B	251	HIS	CA-C-N	7.05	127.03	119.28
1	B	251	HIS	C-N-CA	7.05	127.03	119.28
1	C	4	LEU	N-CA-C	6.94	119.98	109.23
1	C	303	SER	N-CA-C	-6.93	100.22	110.48
1	C	127	SER	N-CA-C	-6.61	105.79	114.31
1	C	330	ALA	N-CA-C	-6.20	103.84	111.40
1	A	119	THR	N-CA-C	-5.99	100.08	109.25
1	A	330	ALA	N-CA-C	-5.92	104.91	111.36
1	A	191	ARG	CA-C-N	-5.87	113.46	122.62
1	A	191	ARG	C-N-CA	-5.87	113.46	122.62
1	B	217	LYS	N-CA-C	-5.83	104.89	112.23
1	A	31	VAL	N-CA-C	-5.79	106.83	111.81
1	B	119	THR	N-CA-C	-5.65	101.49	109.96
1	C	180	SER	N-CA-C	5.54	116.94	108.42
1	B	255	VAL	N-CA-C	-5.42	105.56	110.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	60	ALA	N-CA-C	-5.41	105.49	111.71
1	A	120	ILE	N-CA-C	5.40	115.56	107.51
1	B	226	MET	N-CA-C	-5.40	105.56	111.82
1	B	382	ASP	N-CA-C	-5.39	105.41	111.28
1	A	4	LEU	N-CA-C	5.37	117.24	109.07
1	B	78	VAL	N-CA-C	5.36	116.20	108.48
1	C	187	GLY	N-CA-C	-5.34	107.97	114.92
1	A	78	VAL	N-CA-C	5.33	116.26	108.53
1	A	226	MET	N-CA-C	-5.28	105.69	111.82
1	A	341	ILE	N-CA-C	5.28	115.49	110.42
1	B	256	ILE	N-CA-C	-5.10	102.93	109.30
1	B	4	LEU	N-CA-C	5.04	117.04	109.23

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	3030	0	3053	146	0
1	B	2978	0	3007	144	0
1	C	3030	0	3055	128	0
2	A	23	0	0	0	0
2	B	25	0	0	0	0
2	C	40	0	0	2	0
All	All	9126	0	9115	405	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 22.

All (405) 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:288:ASN:ND2	1:B:289:ARG:H	1.59	1.00
1:C:59:GLU:HA	1:C:62:ALA:HB3	1.46	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:PRO:HG3	1:B:219:SER:OG	1.64	0.97
1:A:208:ARG:H	1:A:208:ARG:HD2	1.30	0.96
1:A:256:ILE:HG12	1:A:271:LEU:HD22	1.52	0.91
1:C:202:THR:HG22	1:C:205:GLN:HG3	1.52	0.90
1:A:124:ASN:HD22	1:A:124:ASN:H	1.18	0.88
1:A:370:ARG:H	1:A:370:ARG:HD2	1.36	0.88
1:C:363:VAL:O	1:C:367:MET:HG2	1.72	0.87
1:B:288:ASN:HD22	1:B:289:ARG:N	1.72	0.86
1:A:391:LYS:HA	1:A:394:MET:HE3	1.57	0.85
1:B:288:ASN:HD22	1:B:289:ARG:H	0.87	0.85
1:A:297:LEU:HD21	1:A:302:LEU:HD21	1.60	0.82
1:A:256:ILE:HG12	1:A:271:LEU:CD2	2.08	0.82
1:C:256:ILE:HG12	1:C:271:LEU:HD22	1.61	0.82
1:A:318:LEU:HD22	1:A:322:MET:HE2	1.63	0.81
1:B:130:THR:HG22	1:B:333:THR:HG23	1.62	0.81
1:B:220:VAL:HG11	1:B:343:VAL:HG13	1.61	0.80
1:A:211:ASN:ND2	1:A:214:MET:HB2	1.96	0.80
1:C:236:ARG:HD2	1:C:241:ALA:O	1.83	0.79
1:A:187:GLY:HA2	1:A:190:PHE:HA	1.65	0.78
1:B:227:ASN:ND2	1:B:228:LYS:H	1.82	0.78
1:A:257:HIS:HD2	1:A:270:GLN:HE22	1.31	0.78
1:B:118:LYS:O	1:B:120:ILE:HD12	1.85	0.76
1:A:278:THR:HB	1:A:279:PRO:HD3	1.68	0.76
1:C:177:GLY:HA2	1:C:263:GLN:NE2	2.01	0.76
1:B:218:ILE:H	1:B:218:ILE:HD12	1.51	0.75
1:B:146:LEU:HB2	1:B:168:LEU:HD12	1.69	0.73
1:B:39:VAL:O	1:B:43:VAL:HG23	1.89	0.73
1:B:52:ARG:HD2	1:B:53:TYR:CE1	2.23	0.73
1:C:52:ARG:HD2	1:C:53:TYR:HE1	1.52	0.73
1:C:52:ARG:HD2	1:C:53:TYR:CE1	2.24	0.73
1:B:394:MET:HE2	1:B:394:MET:HA	1.71	0.72
1:A:204:ASP:OD1	1:A:205:GLN:HG3	1.88	0.72
1:C:181:ILE:HB	1:C:246:MET:CE	2.20	0.72
1:A:257:HIS:CD2	1:A:270:GLN:HE22	2.07	0.72
1:C:41:ARG:HH11	1:C:41:ARG:HG2	1.53	0.72
1:A:120:ILE:HB	1:A:146:LEU:HD23	1.72	0.71
1:A:193:THR:HG23	1:A:196:ARG:H	1.53	0.71
1:A:253:GLN:CD	1:A:253:GLN:H	1.97	0.71
1:A:188:GLY:H	1:A:189:PRO:C	1.98	0.71
1:A:208:ARG:H	1:A:208:ARG:CD	2.01	0.71
1:C:39:VAL:O	1:C:43:VAL:HG23	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:377:ASP:O	1:B:381:VAL:HG23	1.92	0.70
1:C:172:ASP:OD2	1:C:175:GLN:HG2	1.91	0.69
1:B:271:LEU:HD23	1:B:271:LEU:N	2.08	0.69
1:B:299:PHE:HA	1:B:302:LEU:HD22	1.74	0.69
1:B:338:ALA:HB1	1:B:363:VAL:HG21	1.74	0.69
1:C:256:ILE:HG12	1:C:271:LEU:CD2	2.23	0.69
1:C:133:ARG:HG3	1:C:133:ARG:HH11	1.58	0.68
1:A:319:LYS:HA	1:A:322:MET:HE3	1.75	0.68
1:B:148:PRO:HG3	1:B:238:LEU:HD21	1.74	0.68
1:C:318:LEU:HD22	1:C:322:MET:HE3	1.75	0.68
1:A:203:PRO:HD3	1:C:373:GLN:O	1.94	0.68
1:A:12:SER:HB2	1:A:208:ARG:HG3	1.76	0.67
1:A:359:LEU:O	1:A:363:VAL:HG23	1.94	0.67
1:A:370:ARG:H	1:A:370:ARG:CD	2.07	0.67
1:A:319:LYS:O	1:A:323:GLU:HG3	1.95	0.67
1:B:360:ASN:O	1:B:364:LEU:HD23	1.95	0.67
1:A:312:TYR:CE2	1:A:322:MET:HE1	2.30	0.67
1:A:124:ASN:HD22	1:A:124:ASN:N	1.86	0.67
1:C:143:LYS:HA	1:C:143:LYS:HZ3	1.60	0.67
1:C:143:LYS:HA	1:C:143:LYS:NZ	2.10	0.67
1:B:305:LEU:HB2	1:C:305:LEU:HB2	1.75	0.66
1:A:84:ALA:O	1:A:88:MET:HG2	1.95	0.66
1:A:198:LEU:HD22	1:C:111:LEU:HD23	1.79	0.66
1:B:389:ALA:O	1:B:393:VAL:HG23	1.96	0.66
1:B:227:ASN:N	1:B:227:ASN:HD22	1.93	0.65
1:A:390:ARG:HA	1:A:393:VAL:HG12	1.78	0.65
1:A:162:GLN:HB3	1:A:163:PRO:HD3	1.78	0.65
1:A:3:GLN:HB3	1:A:31:VAL:CG2	2.26	0.65
1:A:1:MET:HE3	1:A:29:ARG:HH21	1.62	0.65
1:A:9:SER:HB2	1:A:45:GLN:HE22	1.62	0.65
1:A:196:ARG:HE	1:A:196:ARG:HA	1.61	0.65
1:B:185:GLY:HA3	1:B:228:LYS:HD3	1.79	0.64
1:B:288:ASN:ND2	1:B:289:ARG:N	2.40	0.64
1:B:36:GLY:HA2	1:B:57:ASP:HB2	1.80	0.64
1:C:1:MET:HE3	1:C:29:ARG:NH2	2.11	0.64
1:A:208:ARG:HD2	1:A:208:ARG:N	2.11	0.63
1:B:62:ALA:HB2	1:B:80:SER:HB3	1.79	0.63
1:A:256:ILE:HA	1:A:271:LEU:HD22	1.81	0.63
1:A:185:GLY:HA3	1:A:228:LYS:HE2	1.81	0.62
1:B:17:THR:OG1	1:B:98:MET:HE3	2.00	0.62
1:C:185:GLY:HA3	1:C:228:LYS:HE2	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:LEU:HD12	1:B:45:GLN:NE2	2.13	0.62
1:C:22:ARG:HD3	1:C:48:GLU:OE1	1.99	0.62
1:C:202:THR:HG23	1:C:204:ASP:H	1.64	0.61
1:A:172:ASP:OD1	1:A:175:GLN:HG3	2.00	0.61
1:C:154:ASN:HD21	1:C:282:HIS:CD2	2.17	0.61
1:B:67:THR:O	1:B:71:GLN:HG2	1.99	0.61
1:A:210:PRO:HB3	1:C:115:ARG:HB3	1.82	0.61
1:C:55:VAL:HG22	1:C:79:LEU:HB2	1.82	0.61
1:C:181:ILE:HB	1:C:246:MET:HE1	1.81	0.61
1:A:124:ASN:H	1:A:124:ASN:ND2	1.92	0.61
1:B:162:GLN:CD	1:C:162:GLN:HG3	2.26	0.60
1:C:29:ARG:HH11	1:C:29:ARG:HG2	1.66	0.60
1:A:228:LYS:HA	1:A:228:LYS:HZ2	1.66	0.60
1:B:314:ARG:C	1:B:316:PRO:HD3	2.27	0.60
1:C:181:ILE:HB	1:C:246:MET:HE2	1.83	0.60
1:C:217:LYS:HG3	1:C:343:VAL:HG11	1.85	0.59
1:B:175:GLN:HE21	1:B:175:GLN:HA	1.68	0.59
1:C:165:GLN:NE2	1:C:286:TRP:HE1	2.00	0.59
1:C:227:ASN:OD1	1:C:228:LYS:N	2.35	0.58
1:B:271:LEU:HD11	1:C:259:MET:HE1	1.85	0.58
1:B:251:HIS:HD2	1:B:306:THR:O	1.86	0.58
1:B:251:HIS:HE1	1:B:256:ILE:H	1.51	0.58
1:C:253:GLN:O	1:C:254:SER:HB2	2.02	0.58
1:C:278:THR:HB	1:C:279:PRO:CD	2.34	0.58
1:A:327:GLN:HB2	1:A:331:ALA:CB	2.34	0.58
1:A:151:SER:HB2	1:A:276:MET:HE1	1.86	0.57
1:B:87:ASP:O	1:B:91:LEU:HD13	2.05	0.57
1:B:372:PRO:HB3	1:B:378:VAL:HG22	1.86	0.57
1:A:1:MET:HE3	1:A:29:ARG:NH2	2.19	0.57
1:C:253:GLN:CD	1:C:253:GLN:H	2.12	0.57
1:B:384:ASN:O	1:B:388:VAL:HG23	2.05	0.57
1:B:236:ARG:NH2	1:B:243:ALA:HB2	2.21	0.56
1:B:370:ARG:HG2	1:B:371:GLU:N	2.19	0.56
1:C:136:MET:CE	1:C:169:GLY:HA2	2.35	0.56
1:C:65:LEU:C	1:C:65:LEU:HD13	2.30	0.56
1:A:35:ALA:HB3	1:A:42:MET:HE2	1.87	0.56
1:C:59:GLU:HA	1:C:62:ALA:CB	2.29	0.56
1:C:133:ARG:HG3	1:C:133:ARG:NH1	2.18	0.56
1:C:386:ARG:O	1:C:390:ARG:HG3	2.05	0.56
1:C:177:GLY:HA2	1:C:263:GLN:HE21	1.70	0.56
1:A:312:TYR:HE2	1:A:322:MET:HE1	1.68	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:366:LYS:NZ	1:B:395:ARG:HH22	2.04	0.56
1:C:128:LEU:CD1	1:C:238:LEU:HD13	2.36	0.56
1:C:182:LEU:HD23	1:C:247:GLU:HB2	1.86	0.56
1:A:70:GLN:C	1:A:72:GLN:H	2.13	0.56
1:A:226:MET:HE2	1:A:317:CYS:HB3	1.87	0.56
1:A:387:GLU:HG3	1:A:391:LYS:HE3	1.87	0.56
1:C:154:ASN:HD21	1:C:282:HIS:HD2	1.54	0.56
1:A:253:GLN:CD	1:A:253:GLN:N	2.63	0.56
1:A:225:MET:HE1	1:A:250:ILE:CD1	2.37	0.55
1:B:388:VAL:O	1:B:392:GLU:HG2	2.06	0.55
1:B:125:LYS:HD2	1:B:234:GLU:OE1	2.05	0.55
1:C:128:LEU:HD11	1:C:238:LEU:HD13	1.89	0.55
1:C:257:HIS:HD2	1:C:270:GLN:HE22	1.55	0.55
1:C:173:LEU:HD22	1:C:178:VAL:HG11	1.89	0.55
1:A:202:THR:HB	1:A:203:PRO:HD2	1.88	0.55
1:B:374:CYS:C	1:B:376:ASP:H	2.14	0.55
1:A:67:THR:O	1:A:71:GLN:HG2	2.07	0.55
1:A:337:ALA:HB1	1:A:386:ARG:HG3	1.90	0.54
1:B:190:PHE:CZ	1:B:201:MET:HG2	2.42	0.54
1:B:218:ILE:H	1:B:218:ILE:CD1	2.19	0.54
1:A:259:MET:HG2	1:A:269:ALA:HB2	1.89	0.54
1:A:227:ASN:ND2	1:A:228:LYS:N	2.56	0.54
1:B:218:ILE:HD12	1:B:218:ILE:N	2.21	0.54
1:C:236:ARG:NH2	1:C:325:PHE:CD2	2.76	0.54
1:B:341:ILE:HB	1:B:389:ALA:HB1	1.89	0.54
1:A:256:ILE:CG1	1:A:271:LEU:HD22	2.33	0.54
1:A:370:ARG:HD2	1:A:370:ARG:N	2.14	0.54
1:C:13:ILE:HD13	1:C:123:ALA:HB1	1.90	0.53
1:B:139:VAL:O	1:B:143:LYS:N	2.40	0.53
1:C:134:LEU:HD21	1:C:373:GLN:O	2.09	0.53
1:C:135:PHE:O	1:C:139:VAL:HG23	2.09	0.53
1:C:226:MET:HE1	1:C:321:ALA:HB2	1.90	0.53
1:B:124:ASN:H	1:B:124:ASN:HD22	1.55	0.53
1:C:369:MET:SD	1:C:381:VAL:HA	2.49	0.53
1:A:298:ASP:OD2	1:A:301:LYS:HD3	2.09	0.53
1:B:35:ALA:HB2	1:B:42:MET:HE2	1.90	0.53
1:B:227:ASN:HD22	1:B:228:LYS:H	1.52	0.53
1:B:270:GLN:HE21	1:C:266:SER:HB2	1.74	0.53
1:A:18:LEU:HD22	1:A:49:PHE:CE1	2.44	0.53
1:B:276:MET:HA	1:B:276:MET:HE2	1.90	0.53
1:C:389:ALA:O	1:C:393:VAL:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ASP:O	1:A:91:LEU:HD13	2.09	0.52
1:B:124:ASN:HD22	1:B:124:ASN:N	2.07	0.52
1:B:148:PRO:CG	1:B:238:LEU:HD21	2.39	0.52
1:B:374:CYS:C	1:B:376:ASP:N	2.67	0.52
1:B:386:ARG:O	1:B:390:ARG:NH1	2.42	0.52
1:C:117:GLY:HA2	1:C:144:ALA:HB2	1.89	0.52
1:A:191:ARG:HD3	1:A:191:ARG:C	2.35	0.52
1:B:320:LEU:HD11	1:B:361:LEU:HA	1.90	0.52
1:A:256:ILE:HG12	1:A:271:LEU:HD21	1.92	0.52
1:B:189:PRO:HD2	1:B:223:ALA:HA	1.92	0.52
1:C:37:LYS:O	1:C:39:VAL:HG23	2.09	0.52
1:A:122:LEU:HD22	1:A:124:ASN:ND2	2.24	0.52
1:C:162:GLN:HB2	1:C:163:PRO:HD3	1.92	0.52
1:B:84:ALA:O	1:B:88:MET:HG2	2.09	0.52
1:C:33:LEU:HD12	1:C:51:PRO:HG3	1.92	0.52
1:C:256:ILE:HD13	1:C:259:MET:HE3	1.92	0.52
1:C:311:ASP:HB3	1:C:313:ASP:OD1	2.10	0.52
1:B:175:GLN:HA	1:B:175:GLN:NE2	2.25	0.52
1:B:361:LEU:C	1:B:361:LEU:HD13	2.35	0.52
1:B:64:LEU:O	1:B:68:MET:HG3	2.10	0.51
1:B:236:ARG:HD2	1:B:241:ALA:O	2.11	0.51
1:A:130:THR:HB	1:A:378:VAL:CG1	2.40	0.51
1:B:338:ALA:CB	1:B:363:VAL:HG21	2.39	0.51
1:C:150:ASP:OD2	1:C:153:HIS:ND1	2.44	0.51
1:A:125:LYS:HD3	1:A:150:ASP:HB2	1.92	0.51
1:B:172:ASP:OD2	1:B:175:GLN:HB2	2.11	0.51
1:C:282:HIS:HE1	2:C:423:HOH:O	1.93	0.51
1:A:130:THR:HB	1:A:378:VAL:HG11	1.92	0.51
1:B:36:GLY:CA	1:B:57:ASP:HB2	2.40	0.51
1:C:172:ASP:HB3	1:C:175:GLN:CG	2.41	0.51
1:A:3:GLN:HB3	1:A:31:VAL:HG23	1.93	0.51
1:B:227:ASN:ND2	1:B:228:LYS:N	2.54	0.50
1:B:370:ARG:N	1:B:370:ARG:HD2	2.25	0.50
1:C:172:ASP:HB3	1:C:175:GLN:HG3	1.93	0.50
1:A:154:ASN:O	1:A:158:GLN:HG3	2.12	0.50
1:B:361:LEU:O	1:B:361:LEU:HD22	2.12	0.50
1:A:319:LYS:HA	1:A:322:MET:CE	2.40	0.50
1:B:210:PRO:HA	1:B:219:SER:HB3	1.93	0.50
1:B:36:GLY:O	1:B:56:MET:HE3	2.11	0.50
1:B:196:ARG:NH2	1:B:199:ALA:HB3	2.26	0.50
1:A:133:ARG:HG3	1:A:133:ARG:HH11	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:GLU:CG	1:A:391:LYS:HE3	2.41	0.50
1:B:374:CYS:HB3	1:B:377:ASP:OD2	2.11	0.50
1:B:297:LEU:HD12	1:B:299:PHE:CE1	2.47	0.50
1:B:304:ALA:HB1	1:C:304:ALA:HB1	1.94	0.50
1:C:259:MET:HG2	1:C:269:ALA:HB2	1.94	0.50
1:C:190:PHE:CE2	1:C:201:MET:HE2	2.46	0.49
1:C:124:ASN:HD22	1:C:124:ASN:N	2.10	0.49
1:C:168:LEU:O	1:C:169:GLY:C	2.55	0.49
1:A:188:GLY:H	1:A:190:PHE:N	2.10	0.49
1:B:166:HIS:CD2	1:B:286:TRP:HZ2	2.29	0.49
1:A:270:GLN:C	1:A:271:LEU:HD23	2.36	0.49
1:B:135:PHE:O	1:B:139:VAL:HG23	2.13	0.49
1:B:207:CYS:C	1:B:209:HIS:H	2.21	0.49
1:B:318:LEU:O	1:B:322:MET:HG3	2.12	0.49
1:B:341:ILE:HG13	1:B:389:ALA:HB3	1.93	0.49
1:C:58:ASP:OD2	1:C:60:ALA:HB3	2.13	0.49
1:C:124:ASN:HD22	1:C:124:ASN:H	1.60	0.49
1:A:150:ASP:OD2	1:A:152:GLU:HG2	2.12	0.49
1:A:211:ASN:HD22	1:A:214:MET:HB2	1.71	0.49
1:B:151:SER:HB2	1:B:276:MET:HE1	1.94	0.49
1:A:102:VAL:HA	1:A:126:GLU:OE1	2.12	0.49
1:A:18:LEU:HD11	1:A:45:GLN:NE2	2.27	0.49
1:A:392:GLU:O	1:A:395:ARG:HG2	2.13	0.49
1:B:89:ALA:HB1	1:B:113:ALA:HB2	1.95	0.49
1:B:366:LYS:HZ1	1:B:395:ARG:HH22	1.59	0.49
1:C:160:LEU:HD13	1:C:164:ILE:HG21	1.95	0.49
1:A:18:LEU:HD22	1:A:49:PHE:CD1	2.48	0.48
1:B:251:HIS:CD2	1:B:306:THR:O	2.66	0.48
1:C:259:MET:HG2	1:C:269:ALA:CB	2.43	0.48
1:A:161:PRO:CG	1:A:164:ILE:HD12	2.43	0.48
1:A:229:GLY:O	1:A:232:TYR:HB3	2.14	0.48
1:A:281:ALA:HB2	1:A:292:SER:HB3	1.95	0.48
1:B:337:ALA:O	1:B:340:GLU:HB2	2.13	0.48
1:A:188:GLY:N	1:A:190:PHE:N	2.62	0.48
1:A:209:HIS:ND1	1:A:210:PRO:HD2	2.28	0.48
1:C:130:THR:HG22	1:C:333:THR:HA	1.95	0.48
1:A:125:LYS:HB2	1:A:234:GLU:OE1	2.13	0.48
1:B:167:ASN:HA	1:B:170:TYR:CZ	2.49	0.48
1:B:341:ILE:HG22	1:B:393:VAL:HG21	1.96	0.48
1:A:188:GLY:N	1:A:189:PRO:C	2.70	0.48
1:B:150:ASP:OD1	1:B:153:HIS:ND1	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:ASN:H	1:C:124:ASN:ND2	2.11	0.48
1:B:62:ALA:CB	1:B:80:SER:HB3	2.43	0.48
1:C:153:HIS:HE1	1:C:234:GLU:OE1	1.96	0.47
1:A:52:ARG:HD2	1:A:53:TYR:CE1	2.49	0.47
1:B:337:ALA:HB2	1:B:382:ASP:OD1	2.13	0.47
1:A:216:ARG:O	1:A:220:VAL:HG23	2.15	0.47
1:A:220:VAL:HG11	1:A:343:VAL:HG13	1.97	0.47
1:A:114:ILE:O	1:A:115:ARG:C	2.58	0.47
1:C:202:THR:HG23	1:C:204:ASP:N	2.30	0.47
1:A:189:PRO:O	1:A:191:ARG:N	2.47	0.47
1:B:194:PRO:HG2	1:B:197:ASP:HB2	1.97	0.47
1:C:157:PHE:CD1	1:C:168:LEU:HD11	2.50	0.46
1:A:122:LEU:HD22	1:A:124:ASN:HD21	1.81	0.46
1:A:340:GLU:HG3	1:A:386:ARG:NH2	2.30	0.46
1:C:18:LEU:HD12	1:C:45:GLN:NE2	2.31	0.46
1:C:122:LEU:HD12	1:C:124:ASN:HD21	1.81	0.46
1:B:36:GLY:C	1:B:57:ASP:HB2	2.41	0.46
1:A:216:ARG:NH1	1:C:83:GLN:HE21	2.14	0.46
1:C:41:ARG:HG2	1:C:41:ARG:NH1	2.24	0.46
1:A:145:GLN:NE2	1:A:286:TRP:CH2	2.84	0.46
1:A:162:GLN:HG3	1:A:166:HIS:HD2	1.80	0.46
1:A:173:LEU:HB3	1:A:178:VAL:HB	1.98	0.46
1:C:191:ARG:HD3	2:C:430:HOH:O	2.16	0.46
1:A:108:LEU:HB2	1:A:109:PRO:HD3	1.98	0.46
1:A:352:ARG:HH11	1:A:352:ARG:HB2	1.80	0.46
1:B:271:LEU:HD11	1:C:259:MET:CE	2.46	0.46
1:B:207:CYS:C	1:B:210:PRO:HD2	2.41	0.46
1:A:206:ALA:HA	1:A:208:ARG:CD	2.46	0.46
1:C:335:LEU:HG	1:C:364:LEU:HD11	1.97	0.46
1:B:256:ILE:HA	1:B:271:LEU:HB3	1.99	0.45
1:B:1:MET:HE3	1:B:29:ARG:HH12	1.82	0.45
1:A:190:PHE:N	1:A:190:PHE:CD1	2.84	0.45
1:A:259:MET:HG2	1:A:269:ALA:CB	2.46	0.45
1:C:29:ARG:HH11	1:C:29:ARG:CG	2.29	0.45
1:C:224:THR:O	1:C:225:MET:HB2	2.16	0.45
1:A:227:ASN:ND2	1:A:228:LYS:H	2.13	0.45
1:B:117:GLY:HA2	1:B:144:ALA:HB2	1.98	0.45
1:B:145:GLN:HE22	1:B:166:HIS:HA	1.81	0.45
1:C:318:LEU:HD22	1:C:322:MET:CE	2.46	0.45
1:A:170:TYR:N	1:A:170:TYR:CD1	2.84	0.45
1:A:390:ARG:O	1:A:393:VAL:HG12	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:ALA:O	1:B:386:ARG:C	2.59	0.45
1:A:128:LEU:CD1	1:A:238:LEU:HG	2.46	0.45
1:A:164:ILE:HD13	1:A:173:LEU:HD23	1.99	0.45
1:A:327:GLN:HB2	1:A:331:ALA:HB2	1.98	0.45
1:B:136:MET:CE	1:B:169:GLY:HA3	2.46	0.45
1:B:177:GLY:HA3	1:C:289:ARG:HB2	1.99	0.45
1:B:251:HIS:CE1	1:B:256:ILE:H	2.32	0.45
1:A:225:MET:HE2	1:A:228:LYS:HG3	1.98	0.45
1:A:275:ASP:OD1	1:A:277:ARG:HD3	2.16	0.45
1:B:257:HIS:HD2	1:B:270:GLN:OE1	1.99	0.45
1:C:120:ILE:N	1:C:120:ILE:HD12	2.32	0.45
1:C:285:ALA:O	1:C:286:TRP:C	2.59	0.45
1:C:29:ARG:CZ	1:C:29:ARG:HB2	2.47	0.45
1:C:77:GLU:HG3	1:C:78:VAL:H	1.82	0.45
1:C:226:MET:O	1:C:230:LEU:HG	2.17	0.45
1:B:383:ALA:O	1:B:384:ASN:C	2.60	0.44
1:C:198:LEU:HA	1:C:201:MET:HG2	1.99	0.44
1:B:248:VAL:HG11	1:B:318:LEU:HD11	1.99	0.44
1:C:253:GLN:CD	1:C:253:GLN:N	2.75	0.44
1:B:271:LEU:N	1:B:271:LEU:CD2	2.79	0.44
1:C:36:GLY:O	1:C:56:MET:HE3	2.18	0.44
1:A:352:ARG:NH1	1:A:352:ARG:CB	2.81	0.44
1:B:210:PRO:HD3	1:B:219:SER:CB	2.47	0.44
1:A:310:PRO:HG2	1:A:312:TYR:CZ	2.53	0.44
1:C:125:LYS:C	1:C:127:SER:H	2.24	0.44
1:A:161:PRO:HG2	1:A:164:ILE:HD12	1.99	0.44
1:B:175:GLN:HE21	1:B:175:GLN:CA	2.29	0.44
1:A:65:LEU:HD13	1:A:78:VAL:HG21	1.99	0.44
1:A:66:LYS:O	1:A:70:GLN:HB2	2.18	0.44
1:C:227:ASN:CG	1:C:228:LYS:N	2.75	0.44
1:B:224:THR:HG22	1:B:353:PHE:CZ	2.53	0.43
1:B:367:MET:HE2	1:B:385:ALA:HB2	2.00	0.43
1:C:68:MET:O	1:C:72:GLN:HG2	2.17	0.43
1:B:279:PRO:O	1:B:283:THR:HG23	2.19	0.43
1:B:288:ASN:C	1:B:289:ARG:HG2	2.43	0.43
1:A:146:LEU:HD23	1:A:146:LEU:HA	1.87	0.43
1:A:257:HIS:HD2	1:A:270:GLN:NE2	2.06	0.43
1:A:352:ARG:HB2	1:A:352:ARG:NH1	2.34	0.43
1:C:9:SER:HA	1:C:14:GLY:HA3	2.00	0.43
1:C:108:LEU:HB2	1:C:109:PRO:HD3	2.01	0.43
1:C:132:GLY:O	1:C:136:MET:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LEU:HD12	1:C:259:MET:HB2	2.00	0.43
1:C:352:ARG:HG2	1:C:352:ARG:HH11	1.83	0.43
1:A:65:LEU:HD22	1:A:69:LEU:HG	2.01	0.43
1:A:352:ARG:NE	1:A:354:THR:HG22	2.34	0.43
1:B:356:ILE:HG22	1:B:360:ASN:ND2	2.34	0.43
1:C:229:GLY:O	1:C:232:TYR:HB3	2.17	0.43
1:A:162:GLN:CB	1:A:163:PRO:HD3	2.46	0.43
1:A:352:ARG:CZ	1:A:352:ARG:HB3	2.49	0.43
1:B:270:GLN:HE21	1:C:266:SER:CB	2.32	0.43
1:A:65:LEU:HD13	1:A:78:VAL:CG2	2.49	0.43
1:A:257:HIS:O	1:A:258:SER:CB	2.67	0.42
1:B:363:VAL:O	1:B:367:MET:HG2	2.19	0.42
1:C:257:HIS:CD2	1:C:270:GLN:HE22	2.33	0.42
1:A:256:ILE:HD12	1:A:307:PHE:HZ	1.84	0.42
1:A:297:LEU:HD23	1:A:298:ASP:N	2.33	0.42
1:A:342:THR:O	1:A:351:ILE:HD11	2.19	0.42
1:A:228:LYS:HA	1:A:228:LYS:NZ	2.34	0.42
1:B:205:GLN:C	1:B:207:CYS:H	2.28	0.42
1:B:315:TYR:N	1:B:316:PRO:HD3	2.33	0.42
1:B:345:ALA:O	1:B:350:GLN:HB2	2.19	0.42
1:A:310:PRO:HG2	1:A:312:TYR:CE1	2.54	0.42
1:B:375:VAL:O	1:B:379:LEU:HG	2.19	0.42
1:C:59:GLU:O	1:C:63:LYS:N	2.34	0.42
1:B:35:ALA:HB2	1:B:42:MET:CE	2.49	0.42
1:B:167:ASN:HB2	1:B:171:ALA:HB2	2.01	0.42
1:C:236:ARG:HH11	1:C:242:SER:HA	1.84	0.42
1:B:196:ARG:HH22	1:B:199:ALA:HB3	1.84	0.42
1:B:182:LEU:N	1:B:182:LEU:HD12	2.35	0.42
1:B:356:ILE:O	1:B:360:ASN:ND2	2.52	0.42
1:A:226:MET:HE2	1:A:317:CYS:SG	2.59	0.42
1:B:24:ASN:N	1:B:25:PRO:CD	2.83	0.42
1:B:63:LYS:HE3	1:B:67:THR:OG1	2.19	0.42
1:B:108:LEU:HD23	1:B:108:LEU:HA	1.82	0.42
1:C:136:MET:HE3	1:C:169:GLY:HA2	2.02	0.42
1:C:201:MET:HE3	1:C:205:GLN:HB3	2.00	0.42
1:A:182:LEU:HD11	1:A:261:ARG:HB2	2.03	0.41
1:A:339:ASN:O	1:A:343:VAL:HG23	2.20	0.41
1:B:162:GLN:N	1:B:163:PRO:HD2	2.35	0.41
1:B:201:MET:HG3	1:B:205:GLN:OE1	2.20	0.41
1:C:161:PRO:HD3	1:C:262:TYR:OH	2.19	0.41
1:C:226:MET:HA	1:C:226:MET:HE3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:337:ALA:HB2	1:C:382:ASP:OD1	2.20	0.41
1:B:203:PRO:HB3	1:B:347:LEU:HD23	2.02	0.41
1:C:250:ILE:O	1:C:307:PHE:HA	2.20	0.41
1:A:190:PHE:O	1:A:191:ARG:C	2.63	0.41
1:B:209:HIS:N	1:B:210:PRO:HD2	2.35	0.41
1:B:286:TRP:CD1	1:B:287:PRO:HA	2.55	0.41
1:C:202:THR:HG22	1:C:205:GLN:H	1.84	0.41
1:C:251:HIS:HE1	1:C:305:LEU:HD22	1.85	0.41
1:B:164:ILE:HD13	1:B:171:ALA:HB3	2.02	0.41
1:C:17:THR:OG1	1:C:98:MET:HE3	2.20	0.41
1:A:202:THR:HG22	1:C:373:GLN:HB3	2.02	0.41
1:A:341:ILE:HD13	1:A:386:ARG:HG2	2.03	0.41
1:C:236:ARG:NH2	1:C:325:PHE:HD2	2.18	0.41
1:C:236:ARG:NH1	1:C:242:SER:HA	2.36	0.41
1:A:7:LEU:HA	1:A:34:VAL:HB	2.03	0.41
1:A:224:THR:HG22	1:A:353:PHE:CZ	2.55	0.41
1:A:123:ALA:HA	1:A:149:VAL:HB	2.03	0.41
1:A:190:PHE:H	1:A:190:PHE:HD1	1.61	0.41
1:B:124:ASN:N	1:B:124:ASN:ND2	2.69	0.41
1:B:284:MET:HE3	1:B:284:MET:HB2	1.84	0.41
1:A:18:LEU:CD1	1:A:45:GLN:NE2	2.84	0.41
1:B:199:ALA:HA	1:B:354:THR:HB	2.02	0.41
1:B:345:ALA:HB3	1:B:351:ILE:HD11	2.01	0.41
1:A:98:MET:HE2	1:A:121:LEU:CD1	2.51	0.40
1:B:1:MET:HE3	1:B:29:ARG:NH1	2.36	0.40
1:B:202:THR:OG1	1:B:205:GLN:HG3	2.21	0.40
1:B:210:PRO:HD3	1:B:219:SER:HB3	2.02	0.40
1:A:281:ALA:CB	1:A:292:SER:HB3	2.52	0.40
1:B:314:ARG:HD3	1:B:315:TYR:CZ	2.56	0.40
1:C:122:LEU:CD1	1:C:124:ASN:HD21	2.34	0.40
1:C:348:ALA:HB3	1:C:350:GLN:HG2	2.01	0.40
1:A:36:GLY:O	1:A:56:MET:HE3	2.21	0.40
1:B:329:GLN:HA	1:B:332:THR:OG1	2.22	0.40
1:A:24:ASN:C	1:A:26:GLU:H	2.29	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	396/398 (100%)	367 (93%)	21 (5%)	8 (2%)	6	10
1	B	385/398 (97%)	351 (91%)	29 (8%)	5 (1%)	9	18
1	C	396/398 (100%)	378 (96%)	13 (3%)	5 (1%)	9	18
All	All	1177/1194 (99%)	1096 (93%)	63 (5%)	18 (2%)	8	16

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	PHE
1	A	210	PRO
1	A	258	SER
1	A	349	GLN
1	B	371	GLU
1	A	191	ARG
1	C	38	ASN
1	C	169	GLY
1	C	258	SER
1	A	211	ASN
1	B	197	ASP
1	A	71	GLN
1	B	199	ALA
1	C	126	GLU
1	B	308	ALA
1	B	367	MET
1	C	368	ASP
1	A	189	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	328/328 (100%)	304 (93%)	24 (7%)	13	27
1	B	323/328 (98%)	300 (93%)	23 (7%)	13	29
1	C	328/328 (100%)	304 (93%)	24 (7%)	13	27
All	All	979/984 (100%)	908 (93%)	71 (7%)	13	27

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LEU
1	A	91	LEU
1	A	111	LEU
1	A	122	LEU
1	A	124	ASN
1	A	152	GLU
1	A	183	LEU
1	A	190	PHE
1	A	191	ARG
1	A	198	LEU
1	A	201	MET
1	A	207	CYS
1	A	208	ARG
1	A	227	ASN
1	A	228	LYS
1	A	236	ARG
1	A	253	GLN
1	A	270	GLN
1	A	271	LEU
1	A	290	VAL
1	A	297	LEU
1	A	318	LEU
1	A	335	LEU
1	A	361	LEU
1	B	4	LEU
1	B	5	THR
1	B	63	LYS
1	B	64	LEU
1	B	72	GLN
1	B	91	LEU
1	B	124	ASN

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Mol	Chain	Res	Type
1	B	162	GLN
1	B	168	LEU
1	B	174	GLU
1	B	201	MET
1	B	210	PRO
1	B	227	ASN
1	B	228	LYS
1	B	236	ARG
1	B	258	SER
1	B	271	LEU
1	B	302	LEU
1	B	306	THR
1	B	318	LEU
1	B	320	LEU
1	B	361	LEU
1	B	368	ASP
1	C	4	LEU
1	C	59	GLU
1	C	93	ASP
1	C	124	ASN
1	C	126	GLU
1	C	140	LYS
1	C	141	GLN
1	C	143	LYS
1	C	183	LEU
1	C	195	LEU
1	C	202	THR
1	C	208	ARG
1	C	228	LYS
1	C	238	LEU
1	C	271	LEU
1	C	290	VAL
1	C	297	LEU
1	C	303	SER
1	C	318	LEU
1	C	320	LEU
1	C	335	LEU
1	C	361	LEU
1	C	376	ASP
1	C	394	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	45	GLN
1	A	72	GLN
1	A	124	ASN
1	A	145	GLN
1	A	162	GLN
1	A	166	HIS
1	A	167	ASN
1	A	175	GLN
1	A	205	GLN
1	A	211	ASN
1	A	227	ASN
1	A	257	HIS
1	A	270	GLN
1	A	329	GLN
1	A	373	GLN
1	B	70	GLN
1	B	71	GLN
1	B	72	GLN
1	B	124	ASN
1	B	145	GLN
1	B	162	GLN
1	B	165	GLN
1	B	166	HIS
1	B	175	GLN
1	B	227	ASN
1	B	251	HIS
1	B	257	HIS
1	B	270	GLN
1	B	288	ASN
1	B	291	ASN
1	B	327	GLN
1	B	384	ASN
1	C	45	GLN
1	C	71	GLN
1	C	83	GLN
1	C	124	ASN
1	C	145	GLN
1	C	165	GLN
1	C	175	GLN
1	C	257	HIS
1	C	282	HIS
1	C	291	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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

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	398/398 (100%)	0.53	14 (3%) 47 42	40, 60, 99, 116	0
1	B	391/398 (98%)	0.76	44 (11%) 10 8	35, 64, 121, 144	0
1	C	398/398 (100%)	0.36	12 (3%) 52 48	25, 52, 91, 109	0
All	All	1187/1194 (99%)	0.55	70 (5%) 28 24	25, 58, 107, 144	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	396	LEU	6.9
1	B	199	ALA	5.7
1	B	218	ILE	4.7
1	A	201	MET	4.4
1	C	170	TYR	4.3
1	B	363	VAL	4.2
1	A	189	PRO	4.1
1	A	202	THR	3.8
1	B	344	ALA	3.8
1	A	73	GLY	3.6
1	A	396	LEU	3.6
1	B	198	LEU	3.5
1	B	364	LEU	3.4
1	B	381	VAL	3.2
1	C	169	GLY	3.1
1	B	393	VAL	2.9
1	C	63	LYS	2.9
1	B	207	CYS	2.9
1	B	200	THR	2.8
1	B	376	ASP	2.8
1	B	211	ASN	2.8
1	C	397	ALA	2.8
1	B	375	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	370	ARG	2.8
1	C	67	THR	2.8
1	B	57	ASP	2.7
1	A	319	LYS	2.7
1	B	380	SER	2.6
1	B	347	LEU	2.6
1	C	62	ALA	2.6
1	A	362	SER	2.6
1	B	328	GLY	2.5
1	A	126	GLU	2.5
1	B	334	ALA	2.5
1	C	138	ALA	2.5
1	A	214	MET	2.5
1	A	394	MET	2.5
1	C	36	GLY	2.4
1	B	373	GLN	2.4
1	B	367	MET	2.4
1	B	324	ALA	2.4
1	B	338	ALA	2.4
1	A	190	PHE	2.4
1	C	212	TRP	2.4
1	A	397	ALA	2.3
1	B	368	ASP	2.3
1	C	66	LYS	2.3
1	B	26	GLU	2.3
1	B	12	SER	2.3
1	B	391	LYS	2.3
1	B	15	CYS	2.3
1	B	392	GLU	2.2
1	A	206	ALA	2.2
1	B	384	ASN	2.2
1	B	69	LEU	2.2
1	B	354	THR	2.2
1	B	395	ARG	2.2
1	B	206	ALA	2.2
1	C	1	MET	2.2
1	B	351	ILE	2.2
1	B	190	PHE	2.2
1	B	126	GLU	2.1
1	A	23	HIS	2.1
1	B	201	MET	2.1
1	B	223	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	383	ALA	2.1
1	B	23	HIS	2.1
1	B	102	VAL	2.1
1	B	388	VAL	2.1
1	C	60	ALA	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.