



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:22 AM UTC

PDB ID : 8HGF / pdb_00008hgf
Title : Human PKM2 mutant - C326S
Authors : Chen, Y.-C.; Lin, K.-T.; Cheng, H.-C.
Deposited on : 2022-11-14
Resolution : 3.10 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

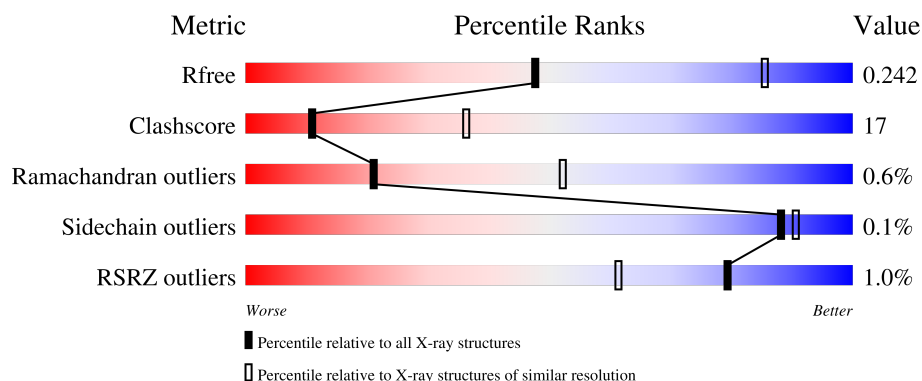
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	551	% 64% 28% • 8%
1	B	551	% 61% 30% • 8%
1	C	551	 56% 18% • 25%
1	D	551	% 62% 28% • 8%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 30019 atoms, of which 15162 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 Pyruvate kinase PKM.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	508	Total	C	H	N	O	S	0	0	0
			7865	2447	3975	689	731	23			
1	B	508	Total	C	H	N	O	S	0	0	0
			7861	2447	3971	689	731	23			
1	D	508	Total	C	H	N	O	S	0	0	0
			7860	2447	3970	689	731	23			
1	C	414	Total	C	H	N	O	S	0	0	0
			6433	2002	3246	575	590	20			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P14618
A	-18	GLY	-	expression tag	UNP P14618
A	-17	SER	-	expression tag	UNP P14618
A	-16	SER	-	expression tag	UNP P14618
A	-15	HIS	-	expression tag	UNP P14618
A	-14	HIS	-	expression tag	UNP P14618
A	-13	HIS	-	expression tag	UNP P14618
A	-12	ARG	-	expression tag	UNP P14618
A	-11	HIS	-	expression tag	UNP P14618
A	-10	HIS	-	expression tag	UNP P14618
A	-9	SER	-	expression tag	UNP P14618
A	-8	SER	-	expression tag	UNP P14618
A	-7	GLY	-	expression tag	UNP P14618
A	-6	LEU	-	expression tag	UNP P14618
A	-5	VAL	-	expression tag	UNP P14618
A	-4	PRO	-	expression tag	UNP P14618
A	-3	ARG	-	expression tag	UNP P14618
A	-2	GLY	-	expression tag	UNP P14618
A	-1	SER	-	expression tag	UNP P14618
A	0	HIS	-	expression tag	UNP P14618
A	326	SER	CYS	engineered mutation	UNP P14618

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	initiating methionine	UNP P14618
B	-18	GLY	-	expression tag	UNP P14618
B	-17	SER	-	expression tag	UNP P14618
B	-16	SER	-	expression tag	UNP P14618
B	-15	HIS	-	expression tag	UNP P14618
B	-14	HIS	-	expression tag	UNP P14618
B	-13	HIS	-	expression tag	UNP P14618
B	-12	ARG	-	expression tag	UNP P14618
B	-11	HIS	-	expression tag	UNP P14618
B	-10	HIS	-	expression tag	UNP P14618
B	-9	SER	-	expression tag	UNP P14618
B	-8	SER	-	expression tag	UNP P14618
B	-7	GLY	-	expression tag	UNP P14618
B	-6	LEU	-	expression tag	UNP P14618
B	-5	VAL	-	expression tag	UNP P14618
B	-4	PRO	-	expression tag	UNP P14618
B	-3	ARG	-	expression tag	UNP P14618
B	-2	GLY	-	expression tag	UNP P14618
B	-1	SER	-	expression tag	UNP P14618
B	0	HIS	-	expression tag	UNP P14618
B	326	SER	CYS	engineered mutation	UNP P14618
D	-19	MET	-	initiating methionine	UNP P14618
D	-18	GLY	-	expression tag	UNP P14618
D	-17	SER	-	expression tag	UNP P14618
D	-16	SER	-	expression tag	UNP P14618
D	-15	HIS	-	expression tag	UNP P14618
D	-14	HIS	-	expression tag	UNP P14618
D	-13	HIS	-	expression tag	UNP P14618
D	-12	ARG	-	expression tag	UNP P14618
D	-11	HIS	-	expression tag	UNP P14618
D	-10	HIS	-	expression tag	UNP P14618
D	-9	SER	-	expression tag	UNP P14618
D	-8	SER	-	expression tag	UNP P14618
D	-7	GLY	-	expression tag	UNP P14618
D	-6	LEU	-	expression tag	UNP P14618
D	-5	VAL	-	expression tag	UNP P14618
D	-4	PRO	-	expression tag	UNP P14618
D	-3	ARG	-	expression tag	UNP P14618
D	-2	GLY	-	expression tag	UNP P14618
D	-1	SER	-	expression tag	UNP P14618
D	0	HIS	-	expression tag	UNP P14618
D	326	SER	CYS	engineered mutation	UNP P14618

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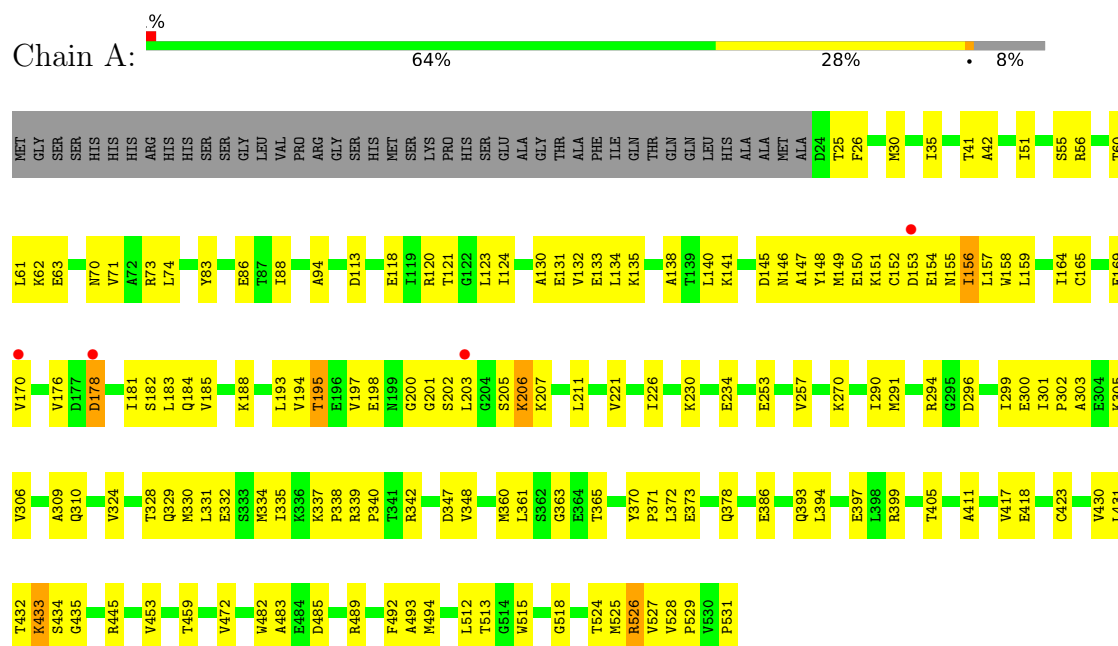
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Chain	Residue	Modelled	Actual	Comment	Reference
C	-19	MET	-	initiating methionine	UNP P14618
C	-18	GLY	-	expression tag	UNP P14618
C	-17	SER	-	expression tag	UNP P14618
C	-16	SER	-	expression tag	UNP P14618
C	-15	HIS	-	expression tag	UNP P14618
C	-14	HIS	-	expression tag	UNP P14618
C	-13	HIS	-	expression tag	UNP P14618
C	-12	ARG	-	expression tag	UNP P14618
C	-11	HIS	-	expression tag	UNP P14618
C	-10	HIS	-	expression tag	UNP P14618
C	-9	SER	-	expression tag	UNP P14618
C	-8	SER	-	expression tag	UNP P14618
C	-7	GLY	-	expression tag	UNP P14618
C	-6	LEU	-	expression tag	UNP P14618
C	-5	VAL	-	expression tag	UNP P14618
C	-4	PRO	-	expression tag	UNP P14618
C	-3	ARG	-	expression tag	UNP P14618
C	-2	GLY	-	expression tag	UNP P14618
C	-1	SER	-	expression tag	UNP P14618
C	0	HIS	-	expression tag	UNP P14618
C	326	SER	CYS	engineered mutation	UNP P14618

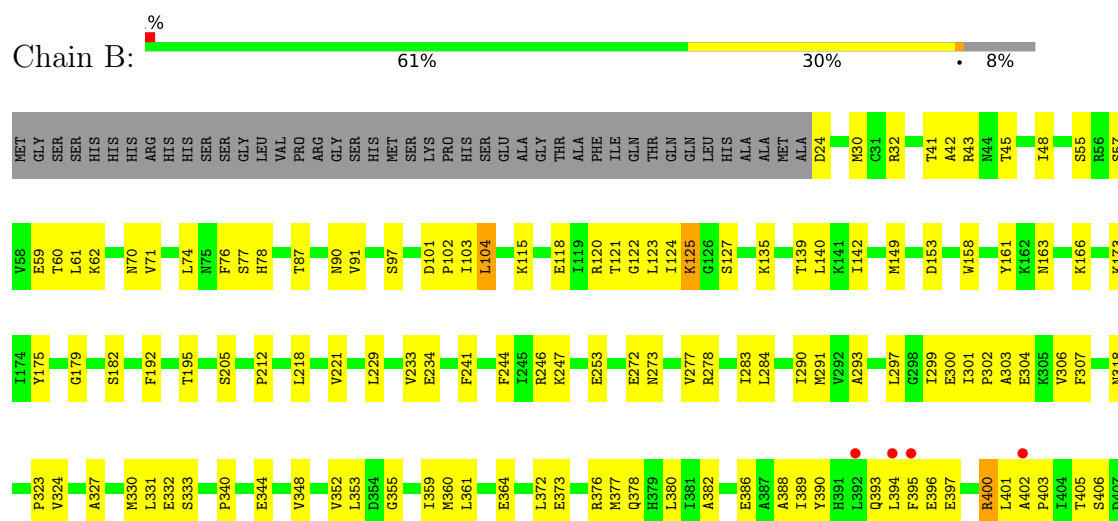
3 Residue-property plots

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: Pyruvate kinase PKM



• Molecule 1: Pyruvate kinase PKM



K498	
G514	
W515	
R516	
P517	
G518	
S519	
M525	
R526	
V527	
P531	

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	181.51Å 156.37Å 121.59Å 90.00° 114.20° 90.00°	Depositor
Resolution (Å)	29.79 – 3.10 29.79 – 3.10	Depositor EDS
% Data completeness (in resolution range)	63.9 (29.79-3.10) 86.2 (29.79-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 3.11Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.182 , 0.237 (Not available) , 0.242	Depositor DCC
R_{free} test set	2000 reflections (4.00%)	wwPDB-VP
Wilson B-factor (Å ²)	53.6	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	30019	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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	0.97	4/3953 (0.1%)	1.08	7/5338 (0.1%)
1	B	0.91	2/3953 (0.1%)	1.04	8/5338 (0.1%)
1	C	0.87	2/3242 (0.1%)	0.97	0/4379
1	D	1.04	4/3953 (0.1%)	1.11	11/5338 (0.2%)
All	All	0.96	12/15101 (0.1%)	1.06	26/20393 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	1
1	C	0	2
1	D	0	2
All	All	0	8

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	432	THR	CA-CB	7.75	1.67	1.53
1	A	434	SER	CB-OG	7.71	1.57	1.42
1	A	178	ASP	CB-CG	6.66	1.68	1.52
1	A	206	LYS	CG-CD	6.03	1.70	1.52
1	C	436	ARG	CG-CD	5.70	1.69	1.52
1	D	402	ALA	C-N	5.69	1.47	1.33
1	D	106	ARG	CG-CD	5.51	1.69	1.52
1	C	423	CYS	CB-SG	-5.47	1.63	1.81
1	B	104	LEU	CG-CD1	5.33	1.70	1.52
1	A	206	LYS	CD-CE	5.30	1.68	1.52
1	D	332	GLU	CG-CD	5.14	1.65	1.52
1	B	432	THR	CB-OG1	5.09	1.51	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	401	LEU	CA-C-N	7.99	141.30	121.80
1	B	401	LEU	C-N-CA	7.99	141.30	121.80
1	D	386	GLU	CA-CB-CG	-6.79	100.52	114.10
1	A	433	LYS	CA-C-N	6.61	129.43	120.38
1	A	433	LYS	C-N-CA	6.61	129.43	120.38
1	B	327	ALA	CA-C-N	-6.56	111.37	122.92
1	B	327	ALA	C-N-CA	-6.56	111.37	122.92
1	D	432	THR	CA-CB-OG1	6.49	119.33	109.60
1	D	401	LEU	CA-C-N	6.26	129.38	120.49
1	D	401	LEU	C-N-CA	6.26	129.38	120.49
1	A	515	TRP	N-CA-C	-6.20	99.14	109.24
1	B	401	LEU	N-CA-C	-6.03	107.16	114.75
1	A	386	GLU	CA-CB-CG	-6.02	102.06	114.10
1	D	432	THR	N-CA-C	-5.90	100.36	109.14
1	D	327	ALA	N-CA-C	5.87	116.02	108.34
1	D	406	SER	N-CA-C	5.84	123.24	110.80
1	B	400	ARG	CB-CG-CD	5.73	124.48	111.30
1	B	32	ARG	CG-CD-NE	-5.70	99.47	112.00
1	D	432	THR	N-CA-CB	5.58	120.66	111.62
1	A	434	SER	CA-CB-OG	5.57	122.23	111.10
1	B	332	GLU	CB-CG-CD	5.38	121.75	112.60
1	D	327	ALA	CA-C-N	5.17	131.41	121.54
1	D	327	ALA	C-N-CA	5.17	131.41	121.54
1	A	526	ARG	CG-CD-NE	-5.08	100.81	112.00
1	A	195	THR	N-CA-C	5.07	116.77	108.76
1	D	127	SER	N-CA-C	-5.05	100.05	110.80

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	423	CYS	Peptide
1	A	518	GLY	Peptide
1	A	526	ARG	Sidechain
1	B	447	ARG	Sidechain
1	C	405	THR	Peptide
1	C	467	ARG	Sidechain
1	D	327	ALA	Peptide
1	D	56	ARG	Sidechain

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	3890	3975	3976	135	1
1	B	3890	3971	3976	162	0
1	C	3187	3246	3249	89	0
1	D	3890	3970	3976	138	1
All	All	14857	15162	15177	496	1

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 17.

All (496) 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:D:118:GLU:OE1	1:D:120:ARG:NH1	1.90	1.04
1:C:71:VAL:HG22	1:C:109:ALA:HB3	1.40	1.03
1:B:400:ARG:HG3	1:C:422:LYS:NZ	1.78	0.98
1:B:395:PHE:CZ	1:B:421:PHE:CE2	2.51	0.98
1:A:41:THR:HG22	1:A:42:ALA:O	1.64	0.98
1:B:400:ARG:HB2	1:C:422:LYS:HE2	1.46	0.96
1:A:140:LEU:HD11	1:A:157:LEU:HB2	1.51	0.93
1:D:402:ALA:HB1	1:D:443:ARG:NH1	1.87	0.89
1:D:56:ARG:NH1	1:D:83:TYR:CE1	2.43	0.87
1:B:330:MET:HE3	1:B:348:VAL:HG22	1.57	0.86
1:C:303:ALA:HA	1:C:306:VAL:HG23	1.57	0.86
1:B:24:ASP:N	1:B:390:TYR:HH	1.76	0.84
1:C:410:GLU:OE1	1:C:440:GLN:NE2	2.11	0.83
1:A:433:LYS:O	1:A:459:THR:HG21	1.80	0.82
1:D:165:CYS:O	1:D:188:LYS:NZ	2.12	0.82
1:A:176:VAL:HB	1:A:181:ILE:HG23	1.61	0.81
1:B:395:PHE:CE1	1:B:421:PHE:CE2	2.69	0.80
1:B:395:PHE:CE1	1:B:421:PHE:CZ	2.72	0.78
1:D:330:MET:HE3	1:D:348:VAL:HG22	1.68	0.76
1:B:400:ARG:HG3	1:C:422:LYS:CE	2.17	0.75
1:B:400:ARG:HG3	1:C:422:LYS:HZ3	1.50	0.75
1:A:165:CYS:O	1:A:188:LYS:NZ	2.18	0.74
1:A:183:LEU:HD13	1:A:195:THR:HG21	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:ARG:NH1	1:D:83:TYR:CZ	2.57	0.73
1:A:121:THR:HG22	1:A:207:LYS:H	1.52	0.73
1:C:405:THR:HG23	1:C:406:SER:HB3	1.71	0.72
1:A:133:GLU:HA	1:A:201:GLY:O	1.90	0.72
1:B:390:TYR:O	1:B:394:LEU:HD13	1.90	0.72
1:A:124:ILE:HA	1:A:152:CYS:HB2	1.70	0.71
1:B:77:SER:HA	1:B:115:LYS:HG3	1.71	0.71
1:B:400:ARG:CB	1:C:422:LYS:HE2	2.21	0.71
1:B:400:ARG:HD2	1:B:403:PRO:HB2	1.73	0.70
1:D:433:LYS:O	1:D:459:THR:HG21	1.91	0.70
1:A:361:LEU:HD21	1:A:378:GLN:NE2	2.07	0.70
1:D:295:GLY:H	1:D:328:THR:HG21	1.55	0.70
1:D:71:VAL:HG22	1:D:109:ALA:HB3	1.74	0.69
1:D:515:TRP:HE1	1:D:520:GLY:HA3	1.57	0.69
1:B:409:THR:HG21	1:B:440:GLN:NE2	2.06	0.69
1:A:361:LEU:HD21	1:A:378:GLN:HE21	1.56	0.69
1:B:518:GLY:O	1:B:519:SER:OG	2.07	0.69
1:D:103:ILE:HA	1:D:496:VAL:HG22	1.73	0.69
1:A:124:ILE:HD11	1:A:130:ALA:C	2.18	0.69
1:B:429:ILE:HD12	1:B:509:VAL:HG21	1.74	0.68
1:D:420:SER:HB3	1:D:428:ILE:HD11	1.74	0.68
1:B:244:PHE:CE2	1:B:246:ARG:HD3	2.29	0.67
1:C:405:THR:OG1	1:C:406:SER:N	2.25	0.67
1:A:197:VAL:O	1:A:197:VAL:HG13	1.92	0.67
1:B:244:PHE:N	1:B:272:GLU:OE2	2.27	0.67
1:A:121:THR:HG22	1:A:207:LYS:N	2.09	0.67
1:A:118:GLU:OE1	1:A:120:ARG:NH1	2.28	0.67
1:C:52:GLY:O	1:C:56:ARG:HG3	1.95	0.67
1:A:124:ILE:HD12	1:A:132:VAL:HG13	1.78	0.66
1:B:405:THR:HG22	1:C:422:LYS:HZ2	1.60	0.66
1:C:397:GLU:O	1:C:401:LEU:HD12	1.95	0.65
1:A:294:ARG:NH2	1:A:347:ASP:OD2	2.30	0.65
1:B:331:LEU:HD23	1:B:344:GLU:HB3	1.78	0.64
1:A:138:ALA:N	1:A:197:VAL:HG12	2.13	0.64
1:A:134:LEU:HD11	1:A:203:LEU:HB2	1.79	0.64
1:D:432:THR:HG23	1:D:435:GLY:H	1.62	0.64
1:B:218:LEU:O	1:B:246:ARG:NH1	2.31	0.63
1:A:348:VAL:HG11	1:A:378:GLN:OE1	1.98	0.63
1:D:361:LEU:HD21	1:D:378:GLN:NE2	2.13	0.63
1:A:56:ARG:HD3	1:A:83:TYR:OH	1.98	0.63
1:A:185:VAL:HA	1:A:195:THR:OG1	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:402:ALA:HB1	1:D:443:ARG:CZ	2.28	0.63
1:D:526:ARG:HD2	1:D:528:VAL:HG12	1.80	0.63
1:D:185:VAL:HA	1:D:195:THR:HG22	1.81	0.62
1:A:334:MET:HA	1:A:337:LYS:O	1.99	0.62
1:D:327:ALA:O	1:D:328:THR:HG22	1.99	0.62
1:B:400:ARG:CG	1:C:422:LYS:HZ3	2.13	0.62
1:A:330:MET:HG3	1:A:360:MET:O	2.00	0.61
1:D:144:LEU:HD11	1:D:164:ILE:HG22	1.82	0.61
1:D:331:LEU:HD23	1:D:334:MET:SD	2.40	0.61
1:D:402:ALA:HB1	1:D:443:ARG:HH12	1.66	0.61
1:D:429:ILE:HD12	1:D:509:VAL:HG21	1.81	0.61
1:B:411:ALA:HA	1:C:422:LYS:CD	2.31	0.61
1:D:103:ILE:HB	1:D:104:LEU:HD12	1.82	0.61
1:A:411:ALA:HB2	1:D:423:CYS:HB3	1.83	0.60
1:D:399:ARG:HE	1:D:400:ARG:HD2	1.66	0.60
1:D:86:GLU:O	1:D:90:ASN:OD1	2.19	0.60
1:C:330:MET:HE3	1:C:348:VAL:HG22	1.82	0.60
1:B:45:THR:HG21	1:B:352:VAL:HG13	1.84	0.60
1:A:124:ILE:HD11	1:A:130:ALA:HB3	1.83	0.60
1:D:125:LYS:HE3	1:D:151:LYS:HD3	1.84	0.59
1:D:168:VAL:HG11	1:D:193:LEU:HD21	1.84	0.59
1:A:124:ILE:HD11	1:A:130:ALA:CB	2.32	0.59
1:A:337:LYS:HG3	1:A:338:PRO:HD2	1.84	0.59
1:B:411:ALA:HA	1:C:422:LYS:HD2	1.84	0.59
1:A:418:GLU:HG3	1:D:418:GLU:HG3	1.85	0.59
1:B:125:LYS:HE2	1:B:153:ASP:HB3	1.85	0.59
1:D:283:ILE:O	1:D:287:SER:OG	2.13	0.59
1:B:57:SER:O	1:B:61:LEU:HD12	2.03	0.58
1:A:290:ILE:O	1:A:324:VAL:HA	2.03	0.58
1:B:101:ASP:OD2	1:B:104:LEU:HD23	2.03	0.58
1:A:489:ARG:O	1:A:492:PHE:HB3	2.04	0.58
1:D:294:ARG:N	1:D:328:THR:HG21	2.19	0.58
1:B:212:PRO:HB3	1:B:299:ILE:HG22	1.85	0.58
1:C:407:ASP:OD1	1:C:409:THR:N	2.35	0.58
1:D:330:MET:CE	1:D:359:ILE:HG23	2.34	0.58
1:A:157:LEU:HD23	1:A:203:LEU:HD11	1.86	0.58
1:C:372:LEU:O	1:C:373:GLU:C	2.48	0.57
1:D:126:GLY:O	1:D:127:SER:OG	2.17	0.57
1:B:24:ASP:N	1:B:390:TYR:OH	2.35	0.57
1:B:429:ILE:CD1	1:B:509:VAL:HG21	2.34	0.57
1:D:192:PHE:O	1:D:193:LEU:HD12	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:LEU:HD22	1:A:513:THR:HG22	1.84	0.57
1:A:123:LEU:O	1:A:124:ILE:HG12	2.04	0.57
1:B:229:LEU:O	1:B:233:VAL:HG23	2.04	0.57
1:A:130:ALA:O	1:A:131:GLU:HG3	2.05	0.57
1:A:157:LEU:HD13	1:A:158:TRP:N	2.20	0.57
1:A:253:GLU:O	1:A:257:VAL:HG23	2.04	0.57
1:D:331:LEU:O	1:D:364:GLU:HG2	2.05	0.56
1:B:393:GLN:HA	1:B:393:GLN:OE1	2.05	0.56
1:D:405:THR:O	1:D:405:THR:HG22	2.05	0.56
1:A:55:SER:HA	1:A:60:THR:HG21	1.87	0.56
1:B:331:LEU:O	1:B:364:GLU:HG2	2.06	0.56
1:D:330:MET:HE2	1:D:359:ILE:HG23	1.87	0.56
1:A:132:VAL:HG21	1:A:153:ASP:O	2.05	0.56
1:B:118:GLU:OE2	1:B:120:ARG:HD3	2.05	0.56
1:D:63:GLU:HB3	1:D:372:LEU:HD21	1.87	0.56
1:D:523:ASN:N	1:D:523:ASN:OD1	2.35	0.56
1:C:301:ILE:HG22	1:C:302:PRO:HD2	1.88	0.56
1:A:56:ARG:NH2	1:A:86:GLU:OE1	2.30	0.56
1:A:294:ARG:HD2	1:A:310:GLN:OE1	2.06	0.56
1:B:175:TYR:CE1	1:B:182:SER:HB3	2.41	0.56
1:B:400:ARG:HD2	1:B:403:PRO:CB	2.36	0.56
1:D:118:GLU:CD	1:D:120:ARG:NH1	2.64	0.56
1:A:169:GLU:O	1:A:185:VAL:HG21	2.06	0.56
1:B:30:MET:HE1	1:B:388:ALA:HB1	1.87	0.56
1:D:420:SER:CB	1:D:428:ILE:HD11	2.36	0.56
1:A:51:ILE:HD12	1:A:61:LEU:HD21	1.88	0.55
1:C:303:ALA:HA	1:C:306:VAL:CG2	2.33	0.55
1:D:26:PHE:CZ	1:D:30:MET:HE3	2.41	0.55
1:A:150:GLU:O	1:A:151:LYS:HG3	2.06	0.55
1:D:109:ALA:HB1	1:D:239:MET:HE3	1.88	0.55
1:C:432:THR:OG1	1:C:435:GLY:HA2	2.07	0.55
1:D:25:THR:HG22	1:D:27:LEU:H	1.70	0.55
1:D:361:LEU:CD2	1:D:378:GLN:NE2	2.69	0.55
1:D:513:THR:OG1	1:D:524:THR:HB	2.07	0.55
1:A:152:CYS:O	1:A:156:ILE:O	2.25	0.55
1:B:400:ARG:CD	1:B:403:PRO:HG2	2.36	0.55
1:D:295:GLY:N	1:D:328:THR:HG21	2.21	0.55
1:A:453:VAL:HG21	1:A:493:ALA:HB2	1.89	0.55
1:B:396:GLU:HG2	1:B:396:GLU:O	2.06	0.55
1:A:296:ASP:O	1:A:300:GLU:HG2	2.07	0.54
1:A:25:THR:HG22	1:A:26:PHE:N	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:LYS:HB3	1:A:194:VAL:HG22	1.90	0.54
1:A:157:LEU:CD2	1:A:203:LEU:HD11	2.37	0.54
1:B:101:ASP:OD1	1:B:102:PRO:HD2	2.08	0.54
1:D:453:VAL:HG21	1:D:493:ALA:HB2	1.88	0.54
1:B:24:ASP:CA	1:B:390:TYR:OH	2.56	0.54
1:A:170:VAL:C	1:A:185:VAL:HG23	2.32	0.54
1:B:59:GLU:O	1:B:59:GLU:HG2	2.06	0.54
1:B:221:VAL:HG21	1:B:253:GLU:CG	2.38	0.54
1:B:453:VAL:HG21	1:B:493:ALA:HB2	1.89	0.54
1:D:302:PRO:HB2	1:D:305:LYS:HD2	1.90	0.54
1:B:515:TRP:HB2	1:B:523:ASN:HB2	1.89	0.53
1:C:327:ALA:HB1	1:C:360:MET:HE2	1.90	0.53
1:B:244:PHE:HE2	1:B:246:ARG:HD3	1.70	0.53
1:A:159:LEU:HD12	1:A:164:ILE:HD13	1.91	0.53
1:B:432:THR:HB	1:B:437:SER:HB2	1.91	0.53
1:D:121:THR:HG22	1:D:122:GLY:O	2.09	0.53
1:B:484:GLU:O	1:B:485:ASP:C	2.52	0.53
1:A:152:CYS:O	1:A:155:ASN:O	2.27	0.53
1:D:392:LEU:O	1:D:396:GLU:OE1	2.27	0.53
1:A:145:ASP:C	1:A:147:ALA:H	2.17	0.53
1:B:432:THR:HB	1:B:437:SER:CB	2.38	0.53
1:C:463:ALA:HB3	1:C:471:PRO:HB3	1.91	0.53
1:B:55:SER:HA	1:B:60:THR:HG21	1.92	0.52
1:D:144:LEU:HD11	1:D:164:ILE:CG2	2.39	0.52
1:D:372:LEU:HD23	1:D:376:ARG:NH1	2.23	0.52
1:A:73:ARG:HD3	1:A:360:MET:SD	2.49	0.52
1:C:323:PRO:HD3	1:C:466:TYR:CE2	2.45	0.52
1:B:59:GLU:O	1:B:59:GLU:CG	2.58	0.52
1:A:135:LYS:O	1:A:197:VAL:HG11	2.09	0.52
1:C:434:SER:HA	1:C:459:THR:HG21	1.91	0.52
1:B:139:THR:HG22	1:B:140:LEU:N	2.23	0.52
1:B:411:ALA:HB2	1:C:422:LYS:HD3	1.92	0.52
1:B:142:ILE:HD12	1:B:195:THR:HG21	1.91	0.52
1:B:330:MET:HG3	1:B:360:MET:O	2.08	0.52
1:D:513:THR:O	1:D:522:THR:HG23	2.10	0.52
1:B:419:ALA:C	1:B:421:PHE:H	2.17	0.52
1:D:304:GLU:HG2	1:D:305:LYS:N	2.24	0.52
1:B:395:PHE:O	1:B:396:GLU:C	2.51	0.52
1:D:187:GLN:HB2	1:D:194:VAL:HB	1.91	0.52
1:C:482:TRP:HH2	1:C:514:GLY:O	1.93	0.52
1:A:159:LEU:HD12	1:A:164:ILE:CD1	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:LEU:HD21	1:B:378:GLN:CG	2.40	0.52
1:A:170:VAL:O	1:A:185:VAL:HG23	2.09	0.51
1:D:399:ARG:HE	1:D:400:ARG:CD	2.22	0.51
1:A:178:ASP:O	1:A:299:ILE:HD11	2.10	0.51
1:B:353:LEU:HD21	1:B:388:ALA:HB3	1.91	0.51
1:B:432:THR:OG1	1:B:437:SER:OG	2.19	0.51
1:A:134:LEU:HD11	1:A:203:LEU:CB	2.39	0.51
1:B:400:ARG:CG	1:C:422:LYS:CE	2.89	0.51
1:D:307:PHE:O	1:D:311:LYS:HG3	2.11	0.51
1:D:399:ARG:NH2	1:D:400:ARG:HH11	2.09	0.51
1:A:154:GLU:HG2	1:A:155:ASN:H	1.75	0.51
1:A:335:ILE:HD11	1:A:363:GLY:HA3	1.92	0.51
1:B:429:ILE:HD12	1:B:509:VAL:CG2	2.40	0.51
1:B:386:GLU:OE1	1:B:467:ARG:NH2	2.44	0.51
1:A:494:MET:HE1	1:A:531:PRO:HD3	1.93	0.50
1:A:303:ALA:C	1:A:305:LYS:H	2.18	0.50
1:B:323:PRO:HB3	1:B:465:LEU:O	2.11	0.50
1:C:115:LYS:HG2	1:C:118:GLU:OE1	2.12	0.50
1:B:103:ILE:HD13	1:B:495:ASN:HB3	1.94	0.50
1:B:498:LYS:HE2	1:B:531:PRO:O	2.12	0.50
1:B:43:ARG:NH2	1:B:70:ASN:OD1	2.44	0.50
1:C:48:ILE:HG12	1:C:71:VAL:HB	1.94	0.50
1:A:176:VAL:HB	1:A:181:ILE:CG2	2.38	0.50
1:A:74:LEU:HD11	1:A:88:ILE:HG13	1.94	0.50
1:B:333:SER:OG	1:B:344:GLU:OE1	2.25	0.50
1:D:74:LEU:HD11	1:D:88:ILE:HG13	1.93	0.50
1:B:330:MET:CE	1:B:348:VAL:HG22	2.36	0.50
1:C:463:ALA:HB1	1:C:469:ILE:HG21	1.93	0.50
1:A:176:VAL:O	1:A:181:ILE:HG22	2.11	0.49
1:D:293:ALA:O	1:D:297:LEU:HB2	2.12	0.49
1:B:470:PHE:N	1:B:470:PHE:CD1	2.80	0.49
1:A:453:VAL:CG2	1:A:493:ALA:HB2	2.42	0.49
1:B:149:MET:HE3	1:B:158:TRP:CH2	2.47	0.49
1:B:241:PHE:CD2	1:B:291:MET:HE1	2.47	0.49
1:B:361:LEU:HD21	1:B:378:GLN:HG3	1.93	0.49
1:B:510:ILE:CD1	1:B:527:VAL:HG22	2.42	0.49
1:D:210:ASN:O	1:D:212:PRO:HD3	2.12	0.49
1:D:297:LEU:O	1:D:301:ILE:HB	2.12	0.49
1:D:123:LEU:O	1:D:152:CYS:HB2	2.13	0.49
1:D:334:MET:HA	1:D:337:LYS:O	2.11	0.49
1:C:321:GLY:CA	1:C:443:ARG:HD2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:ALA:O	1:B:297:LEU:HB2	2.13	0.49
1:C:240:VAL:HG12	1:C:240:VAL:O	2.13	0.49
1:A:405:THR:O	1:D:423:CYS:HA	2.13	0.49
1:C:301:ILE:CG2	1:C:302:PRO:HD2	2.43	0.49
1:C:342:ARG:HD2	1:C:342:ARG:C	2.38	0.49
1:A:405:THR:HG23	1:D:422:LYS:HZ1	1.78	0.48
1:B:406:SER:O	1:B:408:PRO:HD3	2.13	0.48
1:A:331:LEU:HD22	1:A:340:PRO:CB	2.43	0.48
1:B:24:ASP:HA	1:B:390:TYR:OH	2.14	0.48
1:D:253:GLU:O	1:D:257:VAL:HG23	2.14	0.48
1:D:327:ALA:C	1:D:328:THR:HG22	2.38	0.48
1:D:407:ASP:OD1	1:D:407:ASP:O	2.31	0.48
1:C:422:LYS:HG2	1:C:423:CYS:N	2.28	0.48
1:A:56:ARG:HD3	1:A:83:TYR:HH	1.78	0.48
1:A:185:VAL:HG12	1:A:193:LEU:HD21	1.95	0.48
1:B:410:GLU:O	1:B:411:ALA:HB2	2.14	0.48
1:D:67:SER:HB3	1:D:376:ARG:HG3	1.95	0.48
1:B:376:ARG:O	1:B:380:LEU:HD13	2.14	0.48
1:A:140:LEU:O	1:A:195:THR:HG22	2.13	0.48
1:B:456:ASN:O	1:B:457:PRO:C	2.56	0.48
1:D:221:VAL:HG21	1:D:253:GLU:HG3	1.96	0.48
1:B:103:ILE:HG22	1:B:104:LEU:HD22	1.95	0.47
1:B:488:LEU:HD13	1:D:488:LEU:CD1	2.44	0.47
1:B:62:LYS:HE2	1:B:97:SER:CB	2.44	0.47
1:B:359:ILE:HG21	1:B:378:GLN:CD	2.39	0.47
1:A:399:ARG:NH1	1:A:418:GLU:OE2	2.46	0.47
1:B:246:ARG:HH21	1:B:247:LYS:NZ	2.13	0.47
1:B:400:ARG:HD3	1:B:403:PRO:HG2	1.96	0.47
1:B:402:ALA:HB3	1:B:403:PRO:HD3	1.96	0.47
1:A:157:LEU:HD13	1:A:157:LEU:C	2.39	0.47
1:B:74:LEU:HD11	1:B:87:THR:HB	1.96	0.47
1:C:73:ARG:NH1	1:C:113:ASP:OD2	2.47	0.47
1:A:405:THR:HG23	1:D:422:LYS:NZ	2.29	0.47
1:B:233:VAL:O	1:B:234:GLU:C	2.58	0.47
1:D:399:ARG:NE	1:D:400:ARG:HD2	2.30	0.47
1:D:429:ILE:CD1	1:D:509:VAL:HG21	2.45	0.47
1:C:466:TYR:HB2	1:C:469:ILE:HD12	1.96	0.47
1:A:184:GLN:OE1	1:A:198:GLU:OE2	2.33	0.47
1:D:163:ASN:O	1:D:164:ILE:C	2.58	0.47
1:C:30:MET:O	1:C:33:LEU:HG	2.15	0.47
1:A:145:ASP:C	1:A:147:ALA:N	2.73	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:331:LEU:O	1:D:364:GLU:CG	2.63	0.47
1:D:470:PHE:CD2	1:D:502:PHE:HZ	2.33	0.47
1:B:161:TYR:CD1	1:B:161:TYR:C	2.93	0.47
1:D:294:ARG:HD2	1:D:310:GLN:OE1	2.14	0.47
1:A:148:TYR:HB3	1:A:151:LYS:O	2.15	0.47
1:B:173:LYS:HD2	1:B:175:TYR:OH	2.16	0.47
1:D:403:PRO:HD2	1:D:443:ARG:HH22	1.78	0.47
1:D:432:THR:HG21	1:D:437:SER:H	1.79	0.47
1:A:221:VAL:HG12	1:A:226:ILE:HG13	1.98	0.46
1:D:26:PHE:CE1	1:D:30:MET:HE3	2.50	0.46
1:A:399:ARG:NH2	1:D:399:ARG:NH1	2.64	0.46
1:C:321:GLY:HA3	1:C:443:ARG:HD2	1.96	0.46
1:C:324:VAL:O	1:C:357:ASP:HB2	2.15	0.46
1:A:134:LEU:HD12	1:A:134:LEU:H	1.80	0.46
1:A:305:LYS:O	1:A:306:VAL:C	2.58	0.46
1:A:342:ARG:HD3	1:A:342:ARG:C	2.41	0.46
1:B:103:ILE:HD11	1:D:477:PRO:HG3	1.98	0.46
1:B:301:ILE:HG22	1:B:302:PRO:O	2.15	0.46
1:B:395:PHE:CB	1:B:444:TYR:O	2.63	0.46
1:C:58:VAL:O	1:C:59:GLU:C	2.57	0.46
1:A:334:MET:HE3	1:A:339:ARG:HA	1.97	0.46
1:A:365:THR:HA	1:A:371:PRO:HB3	1.98	0.46
1:C:331:LEU:HA	1:C:344:GLU:OE2	2.15	0.46
1:A:329:GLN:HA	1:A:332:GLU:CG	2.46	0.46
1:B:221:VAL:HG21	1:B:253:GLU:HG3	1.98	0.46
1:B:467:ARG:HG2	1:B:468:GLY:N	2.31	0.46
1:D:302:PRO:HG2	1:D:305:LYS:HD2	1.98	0.46
1:A:146:ASN:OD1	1:A:149:MET:HE2	2.15	0.46
1:A:155:ASN:O	1:A:156:ILE:O	2.34	0.46
1:D:407:ASP:OD1	1:D:407:ASP:C	2.58	0.46
1:B:277:VAL:O	1:B:278:ARG:C	2.57	0.46
1:B:303:ALA:HA	1:B:306:VAL:HG23	1.97	0.46
1:B:330:MET:HE2	1:B:359:ILE:HG23	1.98	0.46
1:A:56:ARG:HD3	1:A:83:TYR:CZ	2.51	0.46
1:B:397:GLU:OE1	1:B:397:GLU:HA	2.15	0.46
1:D:431:LEU:HD22	1:D:513:THR:HG22	1.97	0.46
1:D:516:ARG:HB3	1:D:517:PRO:CD	2.46	0.46
1:B:123:LEU:HD23	1:B:205:SER:HB3	1.98	0.46
1:B:359:ILE:HG21	1:B:378:GLN:OE1	2.15	0.46
1:B:400:ARG:CD	1:B:403:PRO:HB2	2.45	0.46
1:D:494:MET:HG2	1:D:531:PRO:HD2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:433:LYS:O	1:B:459:THR:HG21	2.15	0.45
1:B:125:LYS:HE2	1:B:153:ASP:CB	2.47	0.45
1:B:283:ILE:O	1:B:284:LEU:C	2.59	0.45
1:D:158:TRP:CH2	1:D:160:ASP:HB3	2.52	0.45
1:C:372:LEU:O	1:C:375:VAL:N	2.49	0.45
1:C:482:TRP:CH2	1:C:514:GLY:O	2.68	0.45
1:A:205:SER:O	1:A:207:LYS:HE2	2.17	0.45
1:D:507:ASP:O	1:D:530:VAL:HG23	2.16	0.45
1:C:497:GLY:O	1:C:498:LYS:C	2.59	0.45
1:B:163:ASN:HB3	1:B:166:LYS:HD2	1.99	0.45
1:D:290:ILE:CG2	1:D:291:MET:N	2.79	0.45
1:D:123:LEU:HD11	1:D:206:LYS:HE2	1.98	0.45
1:A:62:LYS:HG3	1:A:94:ALA:HB1	1.99	0.45
1:A:331:LEU:HD22	1:A:340:PRO:HB3	1.99	0.45
1:B:142:ILE:CD1	1:B:195:THR:HG21	2.46	0.45
1:B:405:THR:HG22	1:C:422:LYS:NZ	2.30	0.45
1:B:516:ARG:HB3	1:B:517:PRO:CD	2.47	0.45
1:C:372:LEU:HD13	1:C:376:ARG:NH2	2.31	0.45
1:C:470:PHE:CD1	1:C:470:PHE:N	2.84	0.45
1:C:525:MET:HB3	1:C:525:MET:HE2	1.79	0.45
1:A:120:ARG:HD3	1:A:206:LYS:O	2.16	0.45
1:C:372:LEU:C	1:C:374:ALA:N	2.74	0.45
1:C:410:GLU:OE2	1:C:440:GLN:HG2	2.16	0.45
1:A:181:ILE:HD13	1:A:200:GLY:HA2	1.99	0.45
1:B:124:ILE:HD12	1:B:124:ILE:N	2.32	0.45
1:B:421:PHE:HE1	1:B:447:ARG:HD2	1.82	0.45
1:D:96:GLU:O	1:D:97:SER:C	2.60	0.45
1:A:393:GLN:O	1:A:397:GLU:HG3	2.17	0.45
1:D:243:SER:HA	1:D:270:LYS:HB2	1.99	0.44
1:D:426:GLY:O	1:D:427:ALA:HB2	2.16	0.44
1:D:515:TRP:O	1:D:516:ARG:HG2	2.17	0.44
1:B:135:LYS:HB2	1:B:135:LYS:HE2	1.80	0.44
1:B:482:TRP:CH2	1:B:515:TRP:CE3	3.04	0.44
1:B:30:MET:HE1	1:B:388:ALA:CB	2.48	0.44
1:A:472:VAL:CG1	1:A:492:PHE:CE2	3.01	0.44
1:B:372:LEU:HD23	1:B:376:ARG:NH2	2.33	0.44
1:C:334:MET:HA	1:C:337:LYS:O	2.16	0.44
1:A:70:ASN:C	1:A:71:VAL:HG23	2.41	0.44
1:B:175:TYR:CE2	1:B:212:PRO:HG3	2.52	0.44
1:B:400:ARG:CB	1:C:422:LYS:CE	2.93	0.44
1:B:411:ALA:CB	1:C:422:LYS:HD3	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:301:ILE:HG21	1:D:301:ILE:HD13	1.57	0.44
1:C:294:ARG:NH2	1:C:330:MET:SD	2.91	0.44
1:C:48:ILE:HG22	1:C:360:MET:HG3	1.99	0.44
1:C:284:LEU:CD1	1:C:322:LYS:HD2	2.48	0.44
1:C:479:GLN:O	1:C:480:GLU:C	2.60	0.44
1:B:488:LEU:HD13	1:D:488:LEU:HD13	2.00	0.44
1:D:380:LEU:HD12	1:D:380:LEU:HA	1.72	0.44
1:D:489:ARG:O	1:D:492:PHE:HB3	2.18	0.44
1:D:39:PRO:HB2	1:D:383:ARG:HE	1.81	0.44
1:D:164:ILE:HD11	1:D:211:LEU:HD21	1.99	0.44
1:C:103:ILE:HG22	1:C:104:LEU:HD23	2.00	0.44
1:B:175:TYR:HB3	1:B:179:GLY:HA2	1.98	0.43
1:B:340:PRO:HG3	1:B:377:MET:HG2	2.00	0.43
1:B:466:TYR:HB2	1:B:469:ILE:HD12	2.00	0.43
1:B:103:ILE:HG23	1:B:496:VAL:HA	2.00	0.43
1:D:409:THR:HB	1:D:440:GLN:HE22	1.84	0.43
1:C:393:GLN:O	1:C:397:GLU:HG3	2.18	0.43
1:B:48:ILE:HB	1:B:360:MET:HG3	2.00	0.43
1:D:470:PHE:CD1	1:D:470:PHE:N	2.86	0.43
1:C:515:TRP:O	1:C:516:ARG:O	2.36	0.43
1:A:197:VAL:O	1:A:197:VAL:CG1	2.64	0.43
1:B:121:THR:HG22	1:B:122:GLY:N	2.34	0.43
1:B:192:PHE:CD1	1:B:192:PHE:C	2.96	0.43
1:B:432:THR:CB	1:B:437:SER:OG	2.67	0.43
1:B:516:ARG:HB3	1:B:517:PRO:HD2	2.01	0.43
1:D:311:LYS:NZ	1:D:354:ASP:OD1	2.28	0.43
1:A:157:LEU:HD23	1:A:203:LEU:CD1	2.47	0.43
1:B:400:ARG:HD2	1:B:403:PRO:O	2.18	0.43
1:C:406:SER:OG	1:C:407:ASP:N	2.51	0.43
1:A:370:TYR:HB3	1:A:373:GLU:HB2	2.01	0.43
1:D:419:ALA:O	1:D:420:SER:C	2.60	0.43
1:D:109:ALA:HB1	1:D:239:MET:CE	2.49	0.43
1:D:433:LYS:HE3	1:D:455:ARG:H	1.84	0.43
1:D:519:SER:OG	1:D:520:GLY:N	2.49	0.43
1:C:274:HIS:O	1:C:275:GLU:C	2.62	0.43
1:B:395:PHE:HB3	1:B:444:TYR:O	2.18	0.43
1:D:108:VAL:O	1:D:461:ARG:HD2	2.18	0.43
1:A:211:LEU:HD12	1:A:211:LEU:N	2.33	0.43
1:A:230:LYS:HD2	1:A:234:GLU:OE2	2.19	0.43
1:A:430:VAL:HG22	1:A:512:LEU:HB2	2.00	0.43
1:B:103:ILE:C	1:B:104:LEU:HD22	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:290:ILE:HG22	1:D:291:MET:N	2.32	0.43
1:A:51:ILE:HD12	1:A:61:LEU:CD2	2.48	0.43
1:A:63:GLU:HB3	1:A:372:LEU:HD21	2.01	0.43
1:A:329:GLN:HA	1:A:332:GLU:HG2	2.00	0.43
1:B:124:ILE:O	1:B:125:LYS:O	2.37	0.43
1:B:432:THR:HB	1:B:437:SER:OG	2.18	0.43
1:D:112:LEU:C	1:D:112:LEU:HD23	2.44	0.43
1:A:485:ASP:OD2	1:A:489:ARG:NH2	2.52	0.42
1:A:525:MET:HE3	1:A:527:VAL:HG23	2.01	0.42
1:B:139:THR:CG2	1:B:140:LEU:N	2.82	0.42
1:C:45:THR:HG21	1:C:352:VAL:HG22	2.01	0.42
1:C:482:TRP:CD1	1:C:517:PRO:HB3	2.54	0.42
1:A:432:THR:HG21	1:A:435:GLY:HA2	2.00	0.42
1:B:373:GLU:H	1:B:373:GLU:CD	2.26	0.42
1:A:431:LEU:N	1:A:431:LEU:CD1	2.81	0.42
1:B:104:LEU:HD13	1:B:104:LEU:HA	1.96	0.42
1:D:379:HIS:CD2	1:D:379:HIS:C	2.97	0.42
1:A:183:LEU:O	1:A:183:LEU:HD12	2.20	0.42
1:A:472:VAL:HG12	1:A:492:PHE:CE2	2.53	0.42
1:D:395:PHE:O	1:D:396:GLU:C	2.60	0.42
1:C:417:VAL:HG13	1:C:446:PRO:HB3	2.00	0.42
1:B:386:GLU:HA	1:B:389:ILE:HD12	2.00	0.42
1:C:74:LEU:HD11	1:C:88:ILE:HG13	2.01	0.42
1:B:221:VAL:HG21	1:B:253:GLU:HG2	2.01	0.42
1:C:494:MET:HE3	1:C:531:PRO:HD3	2.01	0.42
1:C:515:TRP:C	1:C:516:ARG:O	2.63	0.42
1:A:25:THR:HG22	1:A:26:PHE:H	1.85	0.42
1:B:421:PHE:HE1	1:B:447:ARG:CD	2.32	0.42
1:C:394:LEU:O	1:C:395:PHE:C	2.62	0.42
1:A:394:LEU:HD13	1:A:445:ARG:HG3	2.01	0.42
1:B:90:ASN:O	1:B:91:VAL:C	2.60	0.42
1:D:229:LEU:O	1:D:233:VAL:HG23	2.19	0.42
1:C:515:TRP:O	1:C:516:ARG:HB2	2.18	0.42
1:C:331:LEU:O	1:C:364:GLU:HG2	2.20	0.42
1:C:377:MET:HE2	1:C:377:MET:HB2	1.94	0.42
1:B:382:ALA:O	1:B:386:GLU:HG2	2.20	0.41
1:D:163:ASN:HB3	1:D:166:LYS:CG	2.49	0.41
1:D:361:LEU:HB3	1:D:365:THR:HG23	2.01	0.41
1:B:395:PHE:CZ	1:B:421:PHE:CD2	3.07	0.41
1:A:309:ALA:O	1:A:310:GLN:C	2.62	0.41
1:B:400:ARG:HA	1:B:403:PRO:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:408:PRO:HB3	1:C:527:VAL:HG12	2.02	0.41
1:D:129:THR:O	1:D:131:GLU:N	2.53	0.41
1:A:182:SER:O	1:A:183:LEU:HB3	2.20	0.41
1:B:290:ILE:O	1:B:324:VAL:HA	2.20	0.41
1:B:318:ASN:HD21	1:B:355:GLY:HA3	1.85	0.41
1:B:395:PHE:CZ	1:B:421:PHE:HE2	2.26	0.41
1:D:103:ILE:HG23	1:D:496:VAL:HA	2.02	0.41
1:A:152:CYS:O	1:A:153:ASP:C	2.64	0.41
1:D:302:PRO:HG2	1:D:305:LYS:CD	2.51	0.41
1:C:441:VAL:HG11	1:C:450:ILE:HG12	2.02	0.41
1:A:270:LYS:CG	1:A:291:MET:HE3	2.51	0.41
1:A:301:ILE:O	1:A:302:PRO:C	2.63	0.41
1:B:422:LYS:O	1:C:404:ILE:HB	2.20	0.41
1:B:451:ILE:HD11	1:B:502:PHE:HD2	1.85	0.41
1:D:73:ARG:NH1	1:D:113:ASP:OD2	2.54	0.41
1:D:168:VAL:CG1	1:D:193:LEU:HD21	2.49	0.41
1:C:41:THR:O	1:C:383:ARG:NH2	2.54	0.41
1:C:340:PRO:HG3	1:C:377:MET:HG2	2.02	0.41
1:A:30:MET:HA	1:A:30:MET:HE2	2.03	0.41
1:A:73:ARG:HH11	1:A:113:ASP:CG	2.29	0.41
1:A:512:LEU:HD23	1:A:525:MET:HA	2.03	0.41
1:A:528:VAL:CG2	1:A:529:PRO:HD2	2.50	0.41
1:B:41:THR:HG22	1:B:42:ALA:N	2.35	0.41
1:B:510:ILE:HD13	1:B:527:VAL:HG22	2.01	0.41
1:C:221:VAL:HG22	1:C:253:GLU:OE1	2.21	0.41
1:A:138:ALA:H	1:A:197:VAL:CG1	2.34	0.41
1:A:146:ASN:H	1:A:158:TRP:HE1	1.69	0.41
1:A:303:ALA:C	1:A:305:LYS:N	2.78	0.41
1:B:41:THR:HG22	1:B:42:ALA:H	1.86	0.41
1:B:479:GLN:C	1:B:481:ALA:H	2.29	0.41
1:D:239:MET:HA	1:D:266:LYS:O	2.21	0.41
1:D:297:LEU:HG	1:D:301:ILE:HD12	2.03	0.41
1:A:133:GLU:HA	1:A:202:SER:HA	2.01	0.41
1:A:418:GLU:HG3	1:D:418:GLU:CG	2.51	0.41
1:A:418:GLU:CG	1:D:418:GLU:HG3	2.51	0.41
1:A:512:LEU:HA	1:A:524:THR:O	2.20	0.41
1:B:175:TYR:CD1	1:B:182:SER:HB3	2.55	0.41
1:D:494:MET:CG	1:D:531:PRO:HD2	2.51	0.41
1:C:44:ASN:O	1:C:467:ARG:HG3	2.21	0.41
1:C:219:PRO:O	1:C:220:ALA:HB3	2.20	0.41
1:C:296:ASP:O	1:C:299:ILE:HG12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:VAL:O	1:A:418:GLU:C	2.64	0.40
1:B:48:ILE:HG23	1:B:71:VAL:HB	2.02	0.40
1:B:304:GLU:O	1:B:307:PHE:CE1	2.74	0.40
1:B:423:CYS:HA	1:C:405:THR:HA	2.02	0.40
1:D:248:ALA:O	1:D:249:SER:C	2.64	0.40
1:C:39:PRO:HB2	1:C:383:ARG:HH11	1.86	0.40
1:A:482:TRP:O	1:A:483:ALA:C	2.64	0.40
1:B:70:ASN:HB3	1:B:464:HIS:CG	2.56	0.40
1:D:29:HIS:ND1	1:D:32:ARG:NH2	2.70	0.40
1:D:228:ASP:O	1:D:231:PHE:HB3	2.21	0.40
1:D:512:LEU:C	1:D:513:THR:HG23	2.45	0.40
1:B:76:PHE:C	1:B:78:HIS:N	2.75	0.40
1:B:273:ASN:HB2	1:B:300:GLU:CD	2.45	0.40
1:D:233:VAL:O	1:D:234:GLU:C	2.64	0.40
1:C:517:PRO:C	1:C:519:SER:H	2.29	0.40
1:A:123:LEU:HD12	1:A:150:GLU:HB3	2.04	0.40
1:D:126:GLY:C	1:D:127:SER:HG	2.20	0.40
1:D:330:MET:CE	1:D:359:ILE:CG2	2.99	0.40
1:D:330:MET:HE2	1:D:359:ILE:CG2	2.51	0.40
1:C:372:LEU:O	1:C:374:ALA:N	2.55	0.40
1:D:261:LYS:HD3	1:D:261:LYS:N	2.37	0.40
1:C:39:PRO:HB2	1:C:383:ARG:NH1	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ILE:O	1:D:305:LYS:NZ[2_556]	1.90	0.30

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	506/551 (92%)	456 (90%)	48 (10%)	2 (0%)	30	61
1	B	506/551 (92%)	452 (89%)	50 (10%)	4 (1%)	16	47
1	C	410/551 (74%)	369 (90%)	37 (9%)	4 (1%)	12	41
1	D	506/551 (92%)	456 (90%)	48 (10%)	2 (0%)	30	61
All	All	1928/2204 (88%)	1733 (90%)	183 (10%)	12 (1%)	21	52

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	125	LYS
1	B	411	ALA
1	B	516	ARG
1	C	406	SER
1	A	156	ILE
1	D	127	SER
1	D	519	SER
1	C	328	THR
1	B	127	SER
1	A	328	THR
1	C	403	PRO
1	C	516	ARG

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	419/453 (92%)	419 (100%)	0	100	100
1	B	419/453 (92%)	419 (100%)	0	100	100
1	C	342/453 (76%)	341 (100%)	1 (0%)	86	87
1	D	419/453 (92%)	418 (100%)	1 (0%)	87	89
All	All	1599/1812 (88%)	1597 (100%)	2 (0%)	88	90

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	188	LYS
1	C	422	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	HIS
1	A	210	ASN
1	A	252	HIS
1	A	264	ASN
1	A	329	GLN
1	A	378	GLN
1	B	318	ASN
1	B	440	GLN
1	D	44	ASN
1	D	78	HIS
1	C	252	HIS
1	C	391	HIS

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 ⓘ

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	508/551 (92%)	-0.22	4 (0%) 82 65	25, 45, 74, 89	0
1	B	508/551 (92%)	-0.18	5 (0%) 79 61	31, 51, 76, 98	0
1	C	414/551 (75%)	-0.17	2 (0%) 87 73	31, 50, 72, 94	0
1	D	508/551 (92%)	-0.30	8 (1%) 70 49	27, 46, 73, 96	0
All	All	1938/2204 (87%)	-0.22	19 (0%) 79 61	25, 48, 74, 98	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	402	ALA	4.8
1	C	123	LEU	3.3
1	D	403	PRO	3.2
1	B	395	PHE	3.0
1	D	402	ALA	2.8
1	D	432	THR	2.7
1	D	517	PRO	2.5
1	D	521	PHE	2.4
1	A	170	VAL	2.4
1	D	401	LEU	2.3
1	C	218	LEU	2.3
1	D	518	GLY	2.3
1	A	153	ASP	2.2
1	D	515	TRP	2.2
1	B	521	PHE	2.2
1	B	394	LEU	2.2
1	A	178	ASP	2.2
1	A	203	LEU	2.1
1	B	392	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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.