



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:55 AM UTC

PDB ID : 8GRB / pdb_00008grb
Title : Crystal structure of a constitutively active mutant of the alpha beta heterodimer of human IDH3
Authors : Sun, P.; Chen, X.; Ding, J.
Deposited on : 2022-09-01
Resolution : 2.85 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

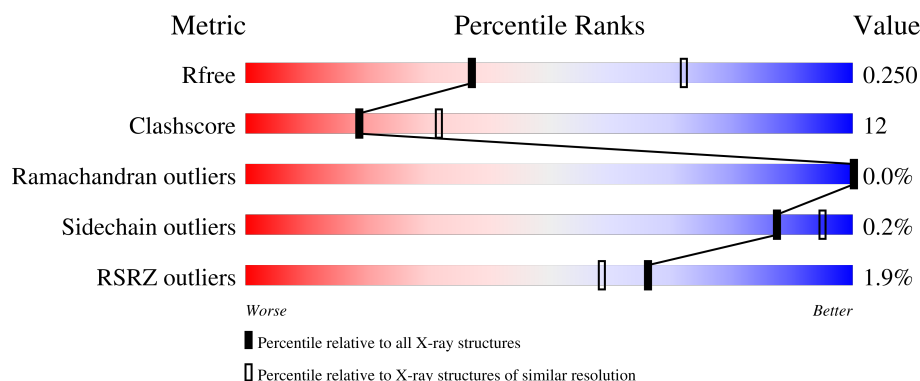
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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.



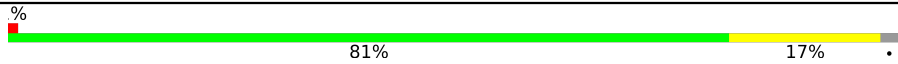
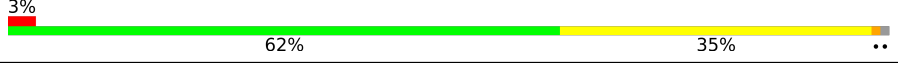
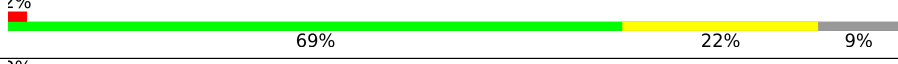
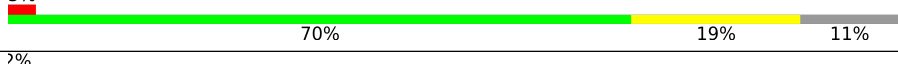
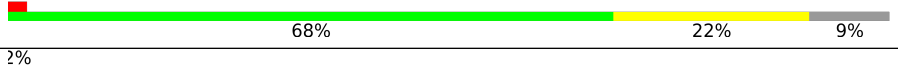
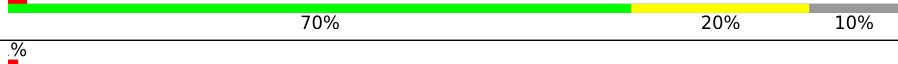
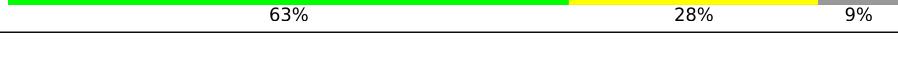
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	339	75% 23% .
1	B	339	% 76% 22% .
1	E	339	% 79% 21% .
1	G	339	% 75% 24% .
1	I	339	% 73% 26% .

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Mol	Chain	Length	Quality of chain
1	K	339	
1	M	339	
1	O	339	
2	C	352	
2	D	352	
2	F	352	
2	H	352	
2	J	352	
2	L	352	
2	N	352	
2	P	352	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 39053 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 Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	333	Total	C	N	O	S	0	0	0
			2474	1553	421	478	22			
1	A	335	Total	C	N	O	S	0	0	0
			2502	1572	428	480	22			
1	E	337	Total	C	N	O	S	0	0	0
			2509	1575	429	483	22			
1	G	335	Total	C	N	O	S	0	0	0
			2483	1561	426	474	22			
1	I	333	Total	C	N	O	S	0	0	0
			2468	1549	421	476	22			
1	K	333	Total	C	N	O	S	0	0	0
			2472	1553	421	476	22			
1	M	335	Total	C	N	O	S	0	0	0
			2490	1565	425	478	22			
1	O	335	Total	C	N	O	S	0	0	0
			2491	1564	426	480	21			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	139	ALA	GLN	engineered mutation	UNP P50213
A	139	ALA	GLN	engineered mutation	UNP P50213
E	139	ALA	GLN	engineered mutation	UNP P50213
G	139	ALA	GLN	engineered mutation	UNP P50213
I	139	ALA	GLN	engineered mutation	UNP P50213
K	139	ALA	GLN	engineered mutation	UNP P50213
M	139	ALA	GLN	engineered mutation	UNP P50213
O	139	ALA	GLN	engineered mutation	UNP P50213

- Molecule 2 is a protein called Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	319	Total	C	N	O	S	0	0	0
			2413	1530	418	445	20			
2	D	319	Total	C	N	O	S	0	0	0
			2391	1509	414	448	20			
2	F	315	Total	C	N	O	S	0	0	0
			2383	1511	414	438	20			
2	H	319	Total	C	N	O	S	0	0	0
			2395	1512	416	447	20			
2	J	317	Total	C	N	O	S	0	0	0
			2406	1525	420	441	20			
2	L	317	Total	C	N	O	S	0	0	0
			2384	1508	416	440	20			
2	N	316	Total	C	N	O	S	0	0	0
			2391	1513	414	444	20			
2	P	320	Total	C	N	O	S	0	0	0
			2401	1516	416	449	20			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	341	GLU	-	expression tag	UNP O43837-2
C	342	ILE	-	expression tag	UNP O43837-2
C	343	CYS	-	expression tag	UNP O43837-2
C	344	ARG	-	expression tag	UNP O43837-2
C	345	ARG	-	expression tag	UNP O43837-2
C	346	VAL	-	expression tag	UNP O43837-2
C	347	LYS	-	expression tag	UNP O43837-2
C	348	ASP	-	expression tag	UNP O43837-2
C	349	LEU	-	expression tag	UNP O43837-2
C	350	ASP	-	expression tag	UNP O43837-2
C	351	GLU	-	expression tag	UNP O43837-2
C	352	ASN	-	expression tag	UNP O43837-2
D	341	GLU	-	expression tag	UNP O43837-2
D	342	ILE	-	expression tag	UNP O43837-2
D	343	CYS	-	expression tag	UNP O43837-2
D	344	ARG	-	expression tag	UNP O43837-2
D	345	ARG	-	expression tag	UNP O43837-2
D	346	VAL	-	expression tag	UNP O43837-2
D	347	LYS	-	expression tag	UNP O43837-2
D	348	ASP	-	expression tag	UNP O43837-2
D	349	LEU	-	expression tag	UNP O43837-2
D	350	ASP	-	expression tag	UNP O43837-2
D	351	GLU	-	expression tag	UNP O43837-2
D	352	ASN	-	expression tag	UNP O43837-2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	341	GLU	-	expression tag	UNP O43837-2
F	342	ILE	-	expression tag	UNP O43837-2
F	343	CYS	-	expression tag	UNP O43837-2
F	344	ARG	-	expression tag	UNP O43837-2
F	345	ARG	-	expression tag	UNP O43837-2
F	346	VAL	-	expression tag	UNP O43837-2
F	347	LYS	-	expression tag	UNP O43837-2
F	348	ASP	-	expression tag	UNP O43837-2
F	349	LEU	-	expression tag	UNP O43837-2
F	350	ASP	-	expression tag	UNP O43837-2
F	351	GLU	-	expression tag	UNP O43837-2
F	352	ASN	-	expression tag	UNP O43837-2
H	341	GLU	-	expression tag	UNP O43837-2
H	342	ILE	-	expression tag	UNP O43837-2
H	343	CYS	-	expression tag	UNP O43837-2
H	344	ARG	-	expression tag	UNP O43837-2
H	345	ARG	-	expression tag	UNP O43837-2
H	346	VAL	-	expression tag	UNP O43837-2
H	347	LYS	-	expression tag	UNP O43837-2
H	348	ASP	-	expression tag	UNP O43837-2
H	349	LEU	-	expression tag	UNP O43837-2
H	350	ASP	-	expression tag	UNP O43837-2
H	351	GLU	-	expression tag	UNP O43837-2
H	352	ASN	-	expression tag	UNP O43837-2
J	341	GLU	-	expression tag	UNP O43837-2
J	342	ILE	-	expression tag	UNP O43837-2
J	343	CYS	-	expression tag	UNP O43837-2
J	344	ARG	-	expression tag	UNP O43837-2
J	345	ARG	-	expression tag	UNP O43837-2
J	346	VAL	-	expression tag	UNP O43837-2
J	347	LYS	-	expression tag	UNP O43837-2
J	348	ASP	-	expression tag	UNP O43837-2
J	349	LEU	-	expression tag	UNP O43837-2
J	350	ASP	-	expression tag	UNP O43837-2
J	351	GLU	-	expression tag	UNP O43837-2
J	352	ASN	-	expression tag	UNP O43837-2
L	341	GLU	-	expression tag	UNP O43837-2
L	342	ILE	-	expression tag	UNP O43837-2
L	343	CYS	-	expression tag	UNP O43837-2
L	344	ARG	-	expression tag	UNP O43837-2
L	345	ARG	-	expression tag	UNP O43837-2
L	346	VAL	-	expression tag	UNP O43837-2

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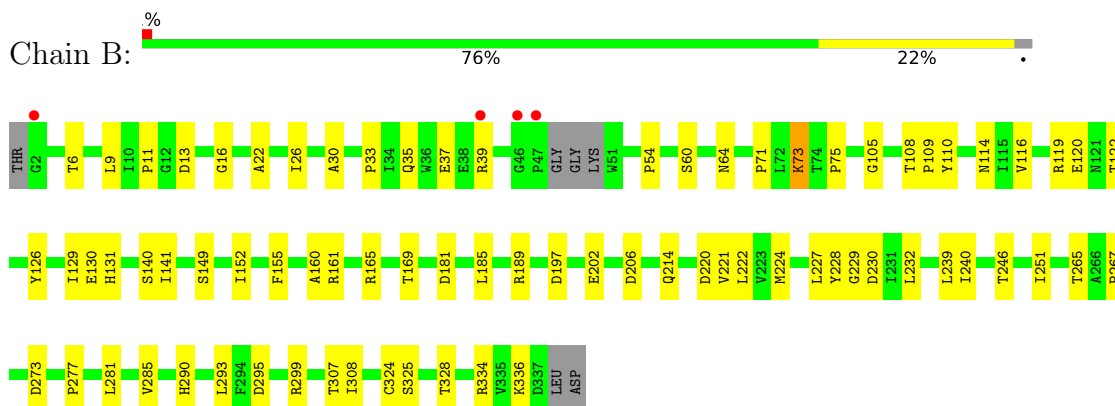
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Chain	Residue	Modelled	Actual	Comment	Reference
L	347	LYS	-	expression tag	UNP O43837-2
L	348	ASP	-	expression tag	UNP O43837-2
L	349	LEU	-	expression tag	UNP O43837-2
L	350	ASP	-	expression tag	UNP O43837-2
L	351	GLU	-	expression tag	UNP O43837-2
L	352	ASN	-	expression tag	UNP O43837-2
N	341	GLU	-	expression tag	UNP O43837-2
N	342	ILE	-	expression tag	UNP O43837-2
N	343	CYS	-	expression tag	UNP O43837-2
N	344	ARG	-	expression tag	UNP O43837-2
N	345	ARG	-	expression tag	UNP O43837-2
N	346	VAL	-	expression tag	UNP O43837-2
N	347	LYS	-	expression tag	UNP O43837-2
N	348	ASP	-	expression tag	UNP O43837-2
N	349	LEU	-	expression tag	UNP O43837-2
N	350	ASP	-	expression tag	UNP O43837-2
N	351	GLU	-	expression tag	UNP O43837-2
N	352	ASN	-	expression tag	UNP O43837-2
P	341	GLU	-	expression tag	UNP O43837-2
P	342	ILE	-	expression tag	UNP O43837-2
P	343	CYS	-	expression tag	UNP O43837-2
P	344	ARG	-	expression tag	UNP O43837-2
P	345	ARG	-	expression tag	UNP O43837-2
P	346	VAL	-	expression tag	UNP O43837-2
P	347	LYS	-	expression tag	UNP O43837-2
P	348	ASP	-	expression tag	UNP O43837-2
P	349	LEU	-	expression tag	UNP O43837-2
P	350	ASP	-	expression tag	UNP O43837-2
P	351	GLU	-	expression tag	UNP O43837-2
P	352	ASN	-	expression tag	UNP O43837-2

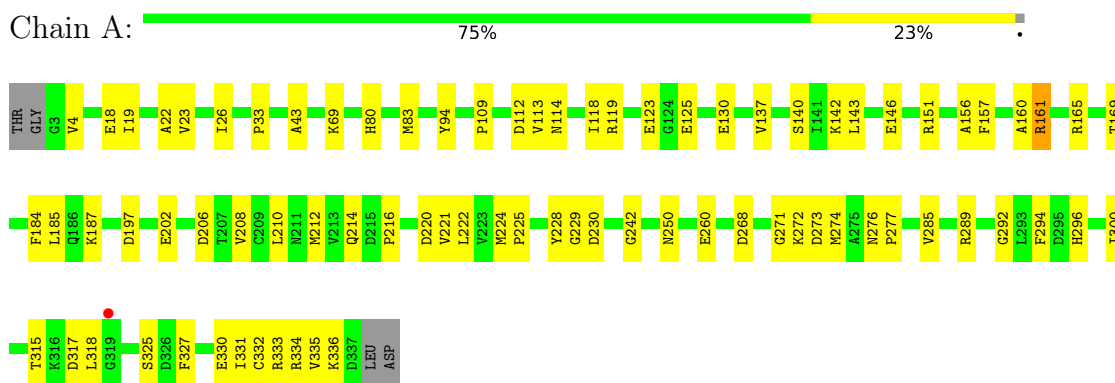
3 Residue-property plots

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.

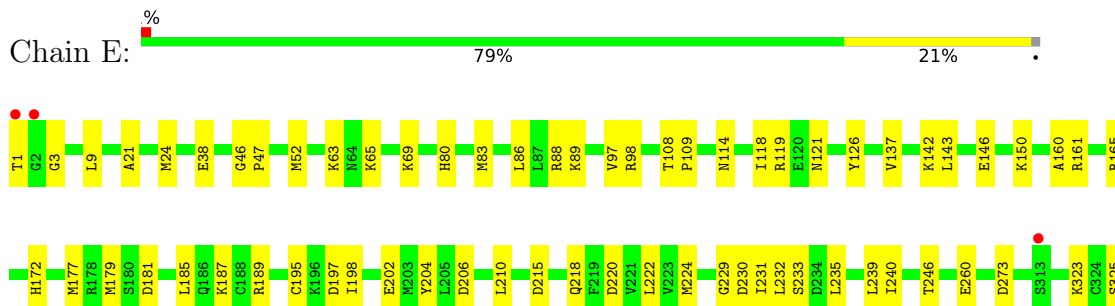
- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

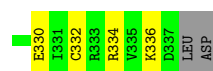


- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

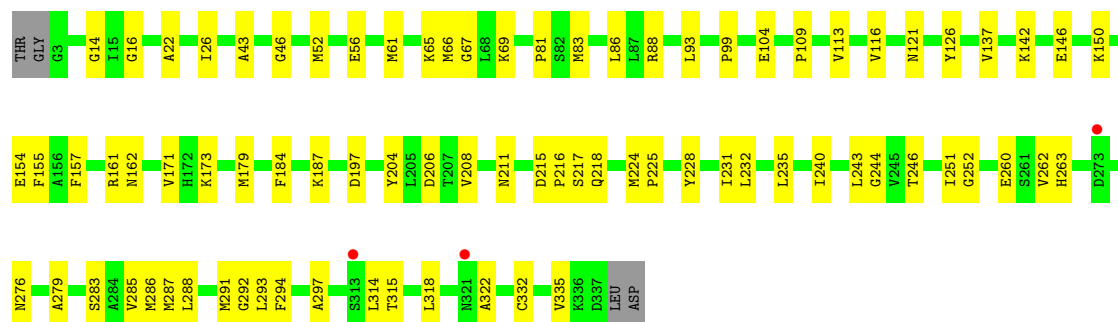
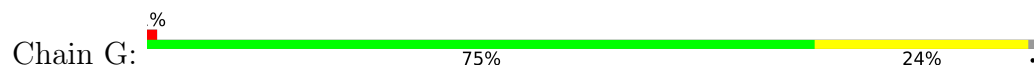


- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

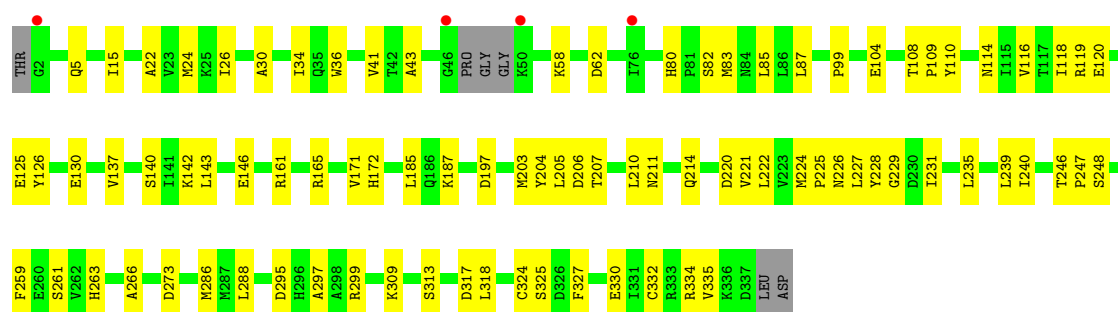




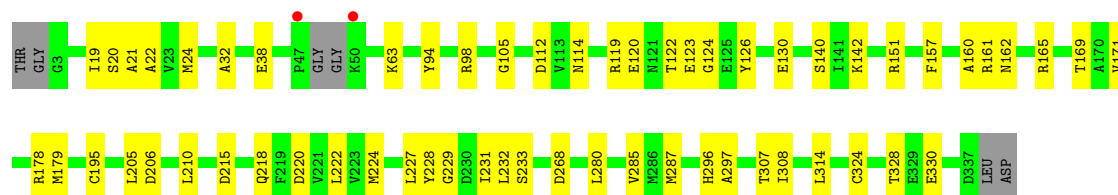
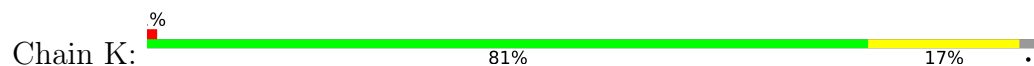
- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial



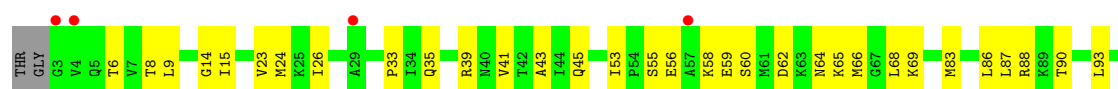
- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

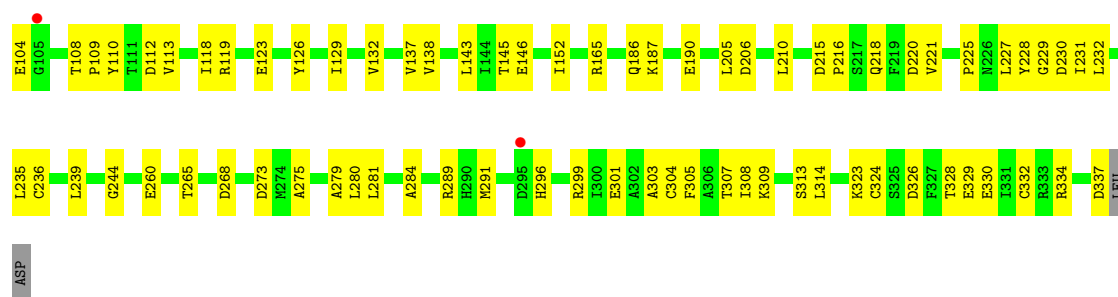


- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

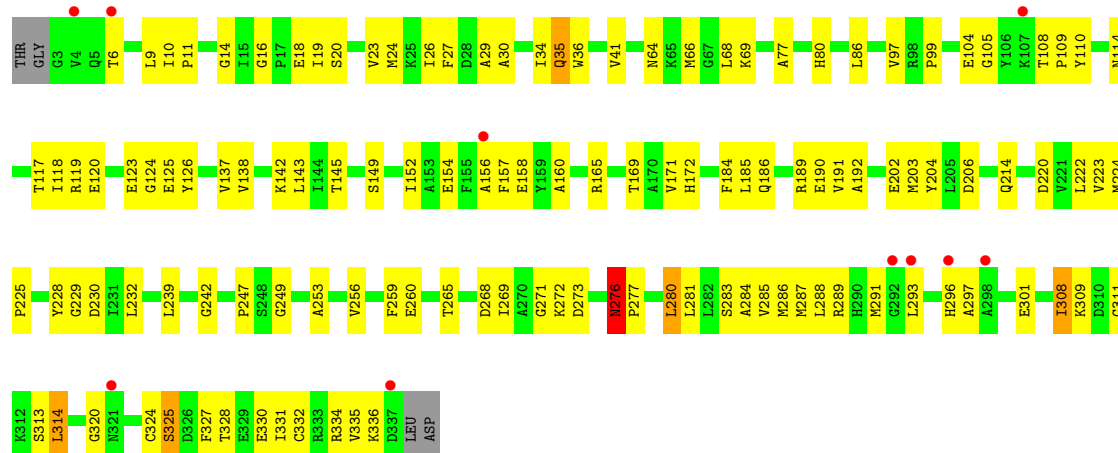


- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

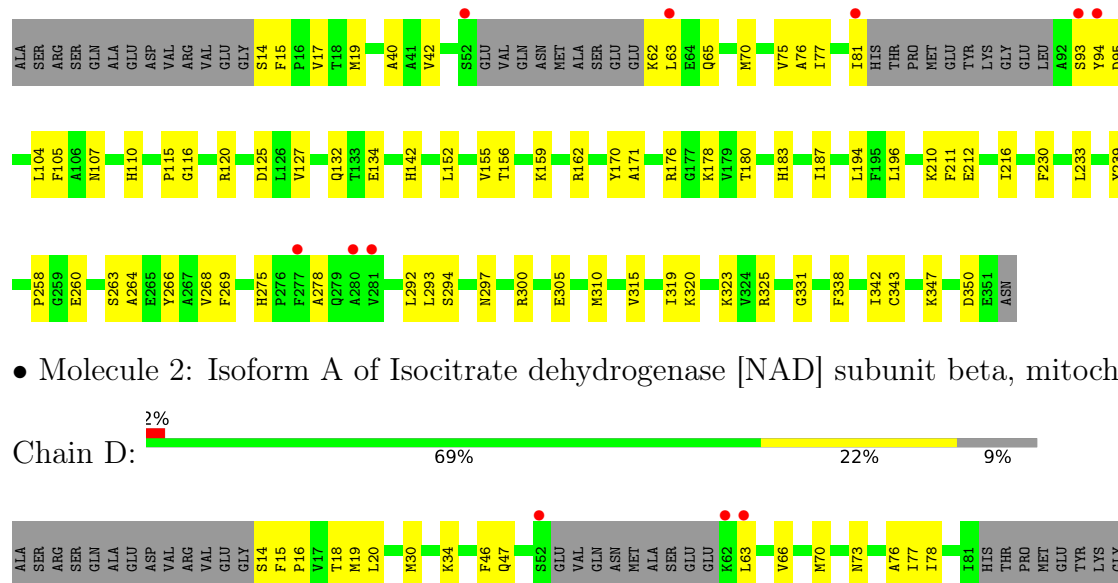




- Molecule 1: Isocitrate dehydrogenase [NAD] subunit alpha, mitochondrial

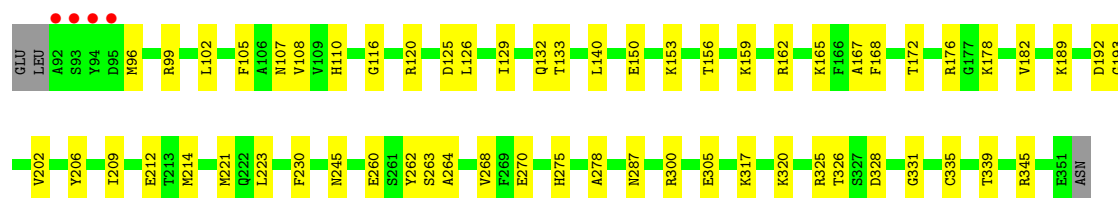


- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial

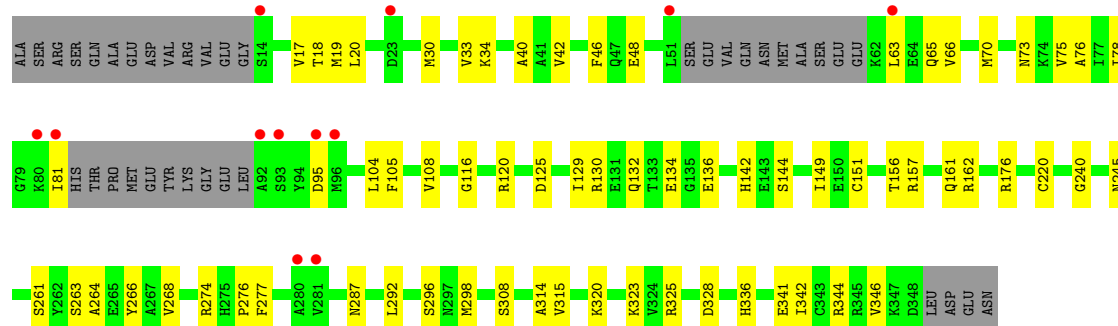


- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial

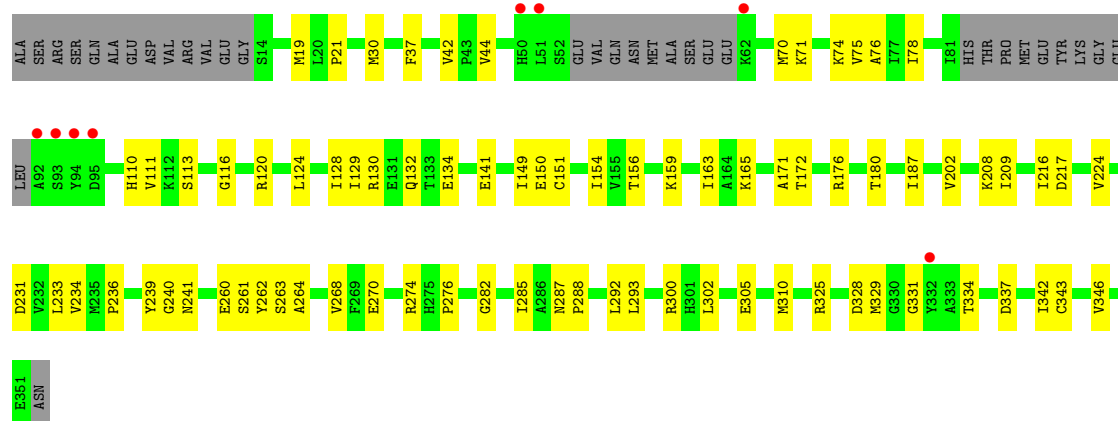




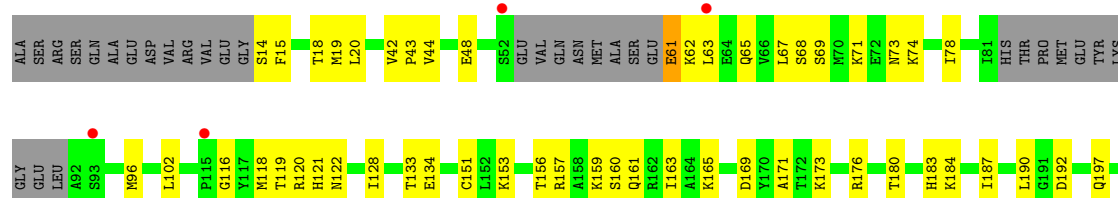
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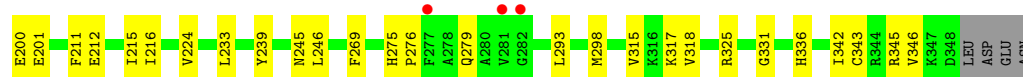


- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial

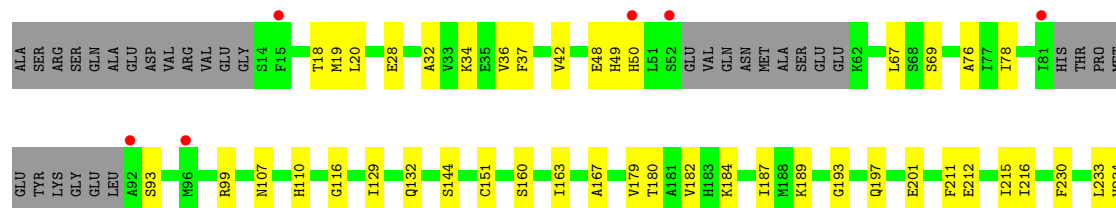


- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial

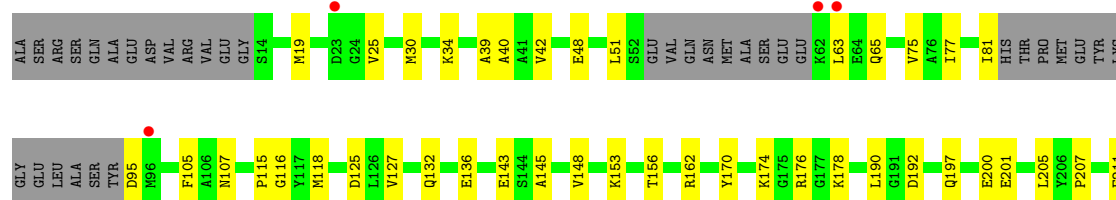




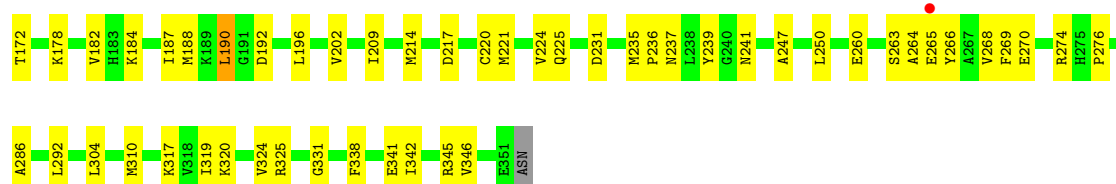
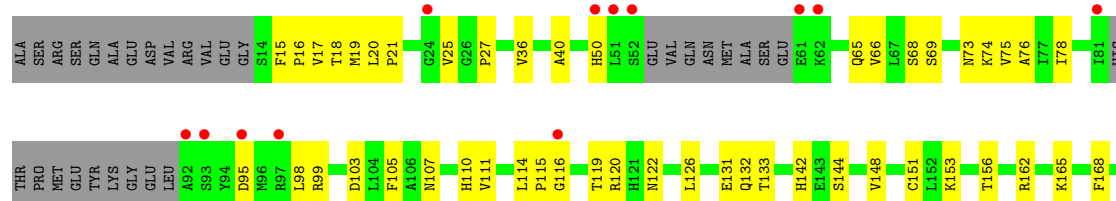
- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial



- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial



- Molecule 2: Isoform A of Isocitrate dehydrogenase [NAD] subunit beta, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	209.15Å 170.03Å 208.69Å 90.00° 103.11° 90.00°	Depositor
Resolution (Å)	49.14 – 2.85 49.14 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.14-2.85) 99.6 (49.14-2.85)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.199 , 0.249 0.200 , 0.250	Depositor DCC
R_{free} test set	8375 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.6	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	0.93	EDS
Total number of atoms	39053	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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	0.40	0/2545	0.59	1/3448 (0.0%)
1	B	0.43	0/2514	0.63	2/3408 (0.1%)
1	E	0.41	0/2552	0.61	0/3459
1	G	0.39	0/2526	0.60	0/3426
1	I	0.45	0/2508	0.64	0/3400
1	K	0.44	0/2512	0.67	0/3406
1	M	0.50	0/2532	0.72	0/3432
1	O	0.57	0/2534	0.83	5/3437 (0.1%)
2	C	0.42	0/2454	0.65	0/3315
2	D	0.41	0/2430	0.64	0/3286
2	F	0.42	0/2424	0.67	3/3275 (0.1%)
2	H	0.39	0/2435	0.61	0/3292
2	J	0.44	0/2447	0.68	1/3304 (0.0%)
2	L	0.42	0/2424	0.67	0/3275
2	N	0.44	0/2431	0.67	0/3286
2	P	0.40	0/2442	0.64	2/3303 (0.1%)
All	All	0.44	0/39710	0.66	14/53752 (0.0%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	73	LYS	CA-CB-CG	6.95	127.99	114.10
1	O	325	SER	CA-CB-OG	6.63	124.36	111.10
2	F	323	LYS	CD-CE-NZ	-6.39	91.46	111.90
1	O	276	ASN	CA-C-N	5.80	126.89	120.45
1	O	276	ASN	C-N-CA	5.80	126.89	120.45
2	F	134	GLU	CB-CA-C	-5.69	107.28	114.40
1	O	308	ILE	N-CA-C	-5.67	105.20	110.53
2	J	63	LEU	N-CA-CB	-5.43	102.09	110.46
1	B	73	LYS	CB-CG-CD	-5.42	98.84	111.30
1	O	26	ILE	CG1-CB-CG2	-5.40	94.51	110.70
2	F	323	LYS	CA-CB-CG	-5.25	103.59	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	190	LEU	CA-C-N	-5.14	111.33	121.41
2	P	190	LEU	C-N-CA	-5.14	111.33	121.41
1	A	161	ARG	NE-CZ-NH1	-5.14	116.36	121.50

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	2502	0	2479	59	0
1	B	2474	0	2429	62	0
1	E	2509	0	2481	57	0
1	G	2483	0	2447	61	0
1	I	2468	0	2419	67	0
1	K	2472	0	2426	52	0
1	M	2490	0	2461	86	0
1	O	2491	0	2450	101	0
2	C	2413	0	2389	52	1
2	D	2391	0	2337	49	0
2	F	2383	0	2361	53	1
2	H	2395	0	2345	57	0
2	J	2406	0	2397	60	0
2	L	2384	0	2352	58	0
2	N	2391	0	2364	56	0
2	P	2401	0	2336	75	0
All	All	39053	0	38473	919	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 12.

All (919) 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)
2:N:51:LEU:HD13	2:N:63:LEU:HD22	1.40	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ARG:NE	1:B:60:SER:HB2	1.79	0.98
1:O:268:ASP:O	1:O:272:LYS:NZ	2.02	0.91
1:M:108:THR:HG21	1:M:239:LEU:O	1.69	0.91
1:K:224:MET:CE	1:K:229:GLY:HA2	2.02	0.89
1:E:330:GLU:OE2	1:E:334:ARG:NH1	2.06	0.87
2:F:76:ALA:HB3	2:F:268:VAL:HG12	1.56	0.87
1:M:330:GLU:HG2	1:M:334:ARG:NH2	1.90	0.86
1:K:161:ARG:NH2	1:K:195:CYS:SG	2.48	0.86
1:I:114:ASN:OD1	1:I:165:ARG:NH2	2.06	0.85
2:L:279:GLN:HE22	2:L:287:ASN:H	1.23	0.85
1:O:332:CYS:O	1:O:335:VAL:HG12	1.78	0.84
2:J:18:THR:HG22	2:J:73:ASN:HD22	1.42	0.83
2:C:81:ILE:HD12	2:C:95:ASP:HA	1.59	0.83
2:N:51:LEU:HD13	2:N:63:LEU:CD2	2.09	0.83
2:F:341:GLU:OE2	2:F:344:ARG:NH2	2.12	0.82
1:E:46:GLY:CA	1:E:52:MET:HE3	2.10	0.82
1:M:6:THR:OG1	1:M:64:ASN:OD1	1.96	0.81
1:A:273:ASP:OD2	1:A:325:SER:OG	1.98	0.81
1:O:23:VAL:HG21	1:O:280:LEU:HD22	1.61	0.81
1:O:34:ILE:HD12	1:O:34:ILE:O	1.81	0.80
1:A:242:GLY:HA2	2:D:221:MET:HE3	1.62	0.80
1:E:1:THR:HG23	1:E:3:GLY:H	1.46	0.80
2:P:18:THR:HG21	2:P:69:SER:HB3	1.63	0.80
1:M:6:THR:HG22	1:M:35:GLN:CG	2.12	0.80
1:A:276:ASN:ND2	1:A:318:LEU:HD13	1.97	0.79
2:N:279:GLN:HE22	2:N:287:ASN:H	1.32	0.78
1:K:224:MET:HE1	1:K:229:GLY:HA2	1.64	0.78
1:A:146:GLU:HB2	1:A:187:LYS:HE3	1.67	0.77
1:M:273:ASP:OD2	1:M:323:LYS:HB2	1.84	0.77
1:E:46:GLY:HA3	1:E:52:MET:HE3	1.67	0.76
1:G:240:ILE:HD11	1:G:246:THR:HG22	1.66	0.76
1:O:137:VAL:HG22	2:P:156:THR:HG22	1.66	0.76
2:J:65:GLN:HA	2:J:68:SER:HB3	1.66	0.76
1:O:169:THR:HB	1:O:222:LEU:HD23	1.67	0.76
1:O:324:CYS:O	1:O:328:THR:OG1	2.03	0.75
2:L:341:GLU:OE1	2:L:341:GLU:O	2.03	0.75
2:J:61:GLU:HG2	2:J:62:LYS:HZ1	1.52	0.75
1:B:265:THR:HG23	1:B:267:PRO:HD3	1.69	0.75
2:L:324:VAL:HG11	2:L:338:PHE:HA	1.69	0.74
1:I:273:ASP:OD2	1:I:325:SER:OG	2.05	0.74
1:I:142:LYS:HB3	2:J:151:CYS:HB3	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:111:VAL:HB	2:P:126:LEU:HB2	1.71	0.73
1:E:114:ASN:OD1	1:E:165:ARG:NH2	2.21	0.73
1:B:39:ARG:HE	1:B:60:SER:HB2	1.51	0.73
1:K:224:MET:HE2	1:K:229:GLY:HA2	1.68	0.73
2:N:51:LEU:CD1	2:N:63:LEU:HD22	2.17	0.72
1:M:108:THR:HG22	1:M:110:TYR:H	1.54	0.72
1:M:330:GLU:HG2	1:M:334:ARG:HH22	1.52	0.72
1:O:114:ASN:OD1	1:O:165:ARG:NH2	2.17	0.72
1:B:37:GLU:HG2	1:B:39:ARG:NH1	2.04	0.72
2:L:182:VAL:HB	2:L:235:MET:HG2	1.71	0.72
2:F:292:LEU:HD13	2:F:342:ILE:HD12	1.71	0.72
2:D:300:ARG:NH1	2:D:305:GLU:OE2	2.22	0.72
1:M:104:GLU:OE2	1:M:309:LYS:HE2	1.90	0.72
1:M:6:THR:HG22	1:M:35:GLN:CD	2.14	0.72
1:K:224:MET:HE2	1:K:229:GLY:CA	2.20	0.71
2:N:30:MET:HE1	2:N:77:ILE:HG12	1.72	0.71
2:C:63:LEU:HD23	2:C:65:GLN:HG3	1.71	0.71
1:I:161:ARG:NH2	1:I:197:ASP:OD1	2.24	0.71
2:F:104:LEU:HD13	2:F:268:VAL:HG21	1.71	0.70
1:I:214:GLN:O	2:J:120:ARG:NH2	2.24	0.70
2:N:81:ILE:HG21	2:N:95:ASP:HB2	1.73	0.70
1:B:240:ILE:HD11	1:B:246:THR:HG22	1.72	0.70
1:G:142:LYS:HB3	2:H:151:CYS:HB3	1.72	0.70
1:I:240:ILE:HD11	1:I:246:THR:HG22	1.72	0.70
1:A:272:LYS:HB3	1:A:274:MET:HG3	1.71	0.70
2:F:34:LYS:NZ	2:F:48:GLU:OE1	2.24	0.70
1:M:33:PRO:HD3	1:M:296:HIS:CE1	2.26	0.70
1:K:308:ILE:HA	1:K:314:LEU:HD21	1.74	0.70
1:G:146:GLU:HG2	1:G:187:LYS:HD2	1.73	0.70
1:E:46:GLY:HA3	1:E:52:MET:CE	2.21	0.69
1:I:118:ILE:CD1	1:I:221:VAL:HG13	2.22	0.69
2:P:21:PRO:HG2	2:P:50:HIS:CD2	2.26	0.69
1:O:214:GLN:O	2:P:120:ARG:NH2	2.24	0.69
2:H:282:GLY:H	2:H:285:ILE:HD12	1.56	0.69
2:P:17:VAL:HG22	2:P:75:VAL:HG23	1.75	0.69
1:K:142:LYS:HB3	2:L:151:CYS:HB3	1.75	0.69
2:J:134:GLU:HG3	2:J:159:LYS:HD2	1.74	0.68
2:C:105:PHE:CE2	2:C:162:ARG:HG2	2.29	0.68
2:P:74:LYS:NZ	2:P:265:GLU:OE2	2.26	0.68
1:B:307:THR:HA	1:B:334:ARG:HH12	1.59	0.68
1:B:108:THR:HG21	1:B:239:LEU:O	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:335:CYS:O	2:L:339:THR:HG23	1.94	0.68
1:O:313:SER:OG	1:O:330:GLU:OE1	2.10	0.68
2:P:20:LEU:HD13	2:P:78:ILE:HG22	1.73	0.68
2:N:335:CYS:O	2:N:339:THR:HG23	1.95	0.67
1:M:6:THR:HA	1:M:35:GLN:O	1.94	0.67
1:G:126:TYR:CD1	2:H:187:ILE:HG21	2.29	0.67
1:I:41:VAL:HG21	1:I:87:LEU:HD21	1.76	0.67
1:I:239:LEU:HD23	2:J:224:VAL:HG11	1.76	0.67
1:G:99:PRO:HB3	1:G:116:VAL:HG22	1.77	0.67
2:D:335:CYS:O	2:D:339:THR:HG22	1.95	0.67
1:E:88:ARG:HD2	1:E:121:ASN:O	1.95	0.67
1:O:165:ARG:NH1	1:O:220:ASP:OD1	2.24	0.67
2:P:110:HIS:HE2	2:P:260:GLU:CD	2.03	0.66
1:G:314:LEU:O	1:G:322:ALA:HB3	1.95	0.66
2:P:76:ALA:HB3	2:P:268:VAL:HG22	1.77	0.66
1:B:105:GLY:HA3	1:B:308:ILE:HG22	1.77	0.66
2:D:76:ALA:HB3	2:D:268:VAL:HG13	1.76	0.66
1:I:259:PHE:HD2	1:I:286:MET:HE3	1.60	0.66
1:E:165:ARG:NH1	1:E:220:ASP:OD1	2.29	0.66
2:C:319:ILE:O	2:C:325:ARG:NH1	2.29	0.66
2:P:236:PRO:HD2	2:P:239:TYR:HD2	1.60	0.65
2:F:116:GLY:HA3	2:F:320:LYS:HA	1.77	0.65
1:O:289:ARG:NH2	1:O:301:GLU:OE2	2.30	0.65
2:C:263:SER:HB3	2:C:266:TYR:H	1.62	0.65
1:E:46:GLY:HA2	1:E:52:MET:HE3	1.78	0.65
2:D:20:LEU:HB2	2:D:78:ILE:HG22	1.78	0.65
1:O:99:PRO:HG2	1:O:247:PRO:HG2	1.78	0.65
1:G:81:PRO:HG2	1:G:86:LEU:HD11	1.79	0.64
1:A:285:VAL:O	1:A:289:ARG:HG3	1.97	0.64
2:D:18:THR:HG22	2:D:73:ASN:HD22	1.63	0.64
2:N:105:PHE:CE2	2:N:162:ARG:HG2	2.31	0.64
1:M:88:ARG:HA	1:M:93:LEU:HD12	1.78	0.64
1:O:171:VAL:HB	1:O:224:MET:HG2	1.79	0.64
2:J:180:THR:HG23	2:J:212:GLU:HB2	1.78	0.64
1:O:330:GLU:OE2	1:O:334:ARG:NH2	2.30	0.63
1:A:185:LEU:HD11	1:A:202:GLU:HG2	1.80	0.63
2:C:110:HIS:NE2	2:C:260:GLU:OE1	2.30	0.63
2:J:18:THR:CG2	2:J:73:ASN:HD22	2.12	0.63
1:O:68:LEU:HD21	1:O:280:LEU:HD21	1.81	0.63
1:B:114:ASN:OD1	1:B:165:ARG:NH2	2.31	0.63
1:A:18:GLU:OE2	1:A:271:GLY:N	2.27	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:287:ASN:HA	2:D:326:THR:HG21	1.81	0.63
1:M:41:VAL:HG21	1:M:87:LEU:HD21	1.79	0.63
2:P:110:HIS:NE2	2:P:260:GLU:OE2	2.29	0.63
1:O:206:ASP:OD1	1:O:206:ASP:N	2.32	0.63
1:O:104:GLU:OE2	1:O:309:LYS:NZ	2.32	0.62
1:M:308:ILE:HA	1:M:314:LEU:HD11	1.81	0.62
2:H:216:ILE:HD13	2:H:239:TYR:HB3	1.81	0.62
2:N:19:MET:HG3	2:N:30:MET:HE3	1.82	0.62
1:B:169:THR:HB	1:B:222:LEU:HD23	1.82	0.62
1:B:39:ARG:NE	1:B:60:SER:CB	2.59	0.62
1:B:129:ILE:HG22	1:B:141:ILE:HG13	1.82	0.62
2:C:125:ASP:OD1	2:C:176:ARG:NH2	2.32	0.62
2:L:197:GLN:NE2	2:L:201:GLU:OE1	2.33	0.62
1:M:6:THR:HG22	1:M:35:GLN:NE2	2.15	0.62
1:M:215:ASP:O	1:M:218:GLN:HG2	2.00	0.62
1:O:142:LYS:HB3	2:P:151:CYS:HB3	1.82	0.62
1:E:146:GLU:HB2	1:E:187:LYS:HD2	1.82	0.61
1:E:179:MET:HG2	2:F:149:ILE:HD12	1.81	0.61
2:L:180:THR:HB	2:L:233:LEU:HD23	1.82	0.61
2:P:114:LEU:HD13	2:P:319:ILE:HD13	1.82	0.61
2:F:261:SER:HB2	2:F:268:VAL:HG23	1.80	0.61
2:H:134:GLU:OE1	2:H:156:THR:OG1	2.06	0.61
1:M:330:GLU:CG	1:M:334:ARG:HH22	2.14	0.61
1:O:19:ILE:O	1:O:23:VAL:HG23	2.01	0.61
1:I:248:SER:C	1:I:286:MET:HE1	2.25	0.61
1:O:18:GLU:OE1	1:O:271:GLY:N	2.33	0.61
2:P:18:THR:CG2	2:P:69:SER:HB3	2.31	0.61
1:B:334:ARG:HG3	1:B:334:ARG:HH11	1.66	0.60
1:I:235:LEU:HD13	2:J:246:LEU:HD11	1.83	0.60
2:C:116:GLY:HA3	2:C:320:LYS:HA	1.83	0.60
1:I:273:ASP:OD1	1:I:324:CYS:HB3	2.01	0.60
1:K:32:ALA:HA	1:K:296:HIS:CD2	2.36	0.60
1:K:160:ALA:HA	1:K:165:ARG:HB2	1.82	0.60
2:D:19:MET:HE3	2:D:46:PHE:CD1	2.36	0.60
1:M:6:THR:CG2	1:M:35:GLN:NE2	2.64	0.60
1:O:27:PHE:HD2	1:O:34:ILE:HD11	1.65	0.60
2:J:20:LEU:HB2	2:J:78:ILE:HG22	1.83	0.60
2:J:153:LYS:NZ	2:J:192:ASP:OD1	2.35	0.60
1:O:30:ALA:HB2	1:O:335:VAL:HG21	1.83	0.60
2:N:281:VAL:HG22	2:N:282:GLY:H	1.67	0.60
1:O:157:PHE:HD2	1:O:191:VAL:HG12	1.67	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:43:ALA:HB2	1:G:83:MET:HE1	1.84	0.60
2:L:37:PHE:CD2	2:L:42:VAL:HG21	2.37	0.60
2:L:48:GLU:OE1	2:L:50:HIS:ND1	2.34	0.60
2:C:110:HIS:HE2	2:C:260:GLU:CD	2.10	0.59
2:P:21:PRO:HB3	2:P:27:PRO:HA	1.83	0.59
1:G:211:ASN:HB3	1:G:218:GLN:OE1	2.02	0.59
2:L:28:GLU:OE2	2:L:284:ASN:N	2.24	0.59
1:M:68:LEU:HD11	1:M:284:ALA:HB2	1.82	0.59
1:M:330:GLU:CG	1:M:334:ARG:NH2	2.64	0.59
2:N:25:VAL:HG21	2:N:286:ALA:HB2	1.84	0.59
1:K:22:ALA:HB1	1:K:328:THR:HG21	1.84	0.59
1:I:207:THR:OG1	2:J:275:HIS:ND1	2.33	0.59
1:B:109:PRO:O	2:C:120:ARG:HD3	2.03	0.59
1:A:161:ARG:NH1	1:A:197:ASP:OD2	2.36	0.59
1:I:24:MET:HG2	1:I:36:TRP:CD1	2.37	0.59
1:M:299:ARG:HH21	1:M:337:ASP:C	2.11	0.59
1:O:160:ALA:HA	1:O:165:ARG:HB2	1.85	0.59
1:M:23:VAL:HG23	1:M:281:LEU:HD12	1.85	0.59
2:D:116:GLY:HA3	2:D:320:LYS:HA	1.84	0.58
1:A:119:ARG:HB2	1:A:229:GLY:HA3	1.83	0.58
2:H:134:GLU:HB3	2:H:154:ILE:O	2.02	0.58
2:C:17:VAL:HG13	2:C:75:VAL:HG23	1.85	0.58
1:I:295:ASP:O	1:I:299:ARG:HG3	2.03	0.58
1:M:146:GLU:HB2	1:M:187:LYS:HD2	1.84	0.58
2:N:39:ALA:HA	2:N:347:LYS:HE3	1.85	0.58
2:F:18:THR:CG2	2:F:73:ASN:HD22	2.16	0.58
2:D:110:HIS:HE2	2:D:260:GLU:CD	2.10	0.58
1:M:9:LEU:HD23	1:M:24:MET:HE3	1.85	0.58
1:B:108:THR:HG22	1:B:110:TYR:H	1.69	0.58
1:G:231:ILE:HA	2:H:217:ASP:HB3	1.86	0.58
2:C:260:GLU:HG2	2:C:269:PHE:CE2	2.39	0.57
1:B:160:ALA:HA	1:B:165:ARG:HB2	1.85	0.57
2:L:341:GLU:OE2	2:L:345:ARG:NH2	2.37	0.57
2:L:99:ARG:HD2	2:L:132:GLN:NE2	2.19	0.57
2:C:300:ARG:NH1	2:C:305:GLU:CB	2.66	0.57
2:F:108:VAL:HG22	2:F:129:ILE:HD12	1.85	0.57
2:H:176:ARG:HD2	2:H:231:ASP:OD1	2.05	0.57
2:P:325:ARG:HB2	2:P:331:GLY:HA3	1.85	0.57
2:H:116:GLY:O	2:H:325:ARG:NH2	2.38	0.57
1:M:289:ARG:NH2	1:M:301:GLU:OE2	2.35	0.57
1:B:73:LYS:O	1:B:75:PRO:HD3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:258:PRO:HB2	2:C:294:SER:HB3	1.87	0.57
1:G:113:VAL:HG22	1:G:216:PRO:HB2	1.86	0.57
1:M:323:LYS:HG3	1:M:326:ASP:CG	2.29	0.57
1:A:169:THR:HB	1:A:222:LEU:HD23	1.86	0.57
1:E:240:ILE:HD11	1:E:246:THR:HG22	1.87	0.57
1:K:215:ASP:O	1:K:218:GLN:HG2	2.05	0.56
1:K:224:MET:CE	1:K:229:GLY:CA	2.76	0.56
2:F:132:GLN:OE1	2:F:274:ARG:NH2	2.36	0.56
1:K:169:THR:HB	1:K:222:LEU:HD23	1.87	0.56
2:L:315:VAL:HG22	2:L:342:ILE:HD13	1.86	0.56
1:B:22:ALA:HB1	1:B:328:THR:HG21	1.87	0.56
2:F:292:LEU:CD1	2:F:342:ILE:HD12	2.36	0.56
2:L:129:ILE:HD12	2:L:167:ALA:HA	1.88	0.56
2:N:178:LYS:HD2	2:N:212:GLU:OE1	2.04	0.56
2:D:18:THR:OG1	2:D:70:MET:HE1	2.05	0.56
1:O:285:VAL:HG13	1:O:297:ALA:HB1	1.87	0.56
1:A:225:PRO:HD2	1:A:228:TYR:HD2	1.71	0.56
1:G:179:MET:HB2	2:H:141:GLU:CD	2.30	0.56
1:I:210:LEU:HB2	2:J:245:ASN:HB3	1.87	0.56
2:N:127:VAL:HB	2:N:232:VAL:HG22	1.85	0.56
2:N:200:GLU:HG3	2:N:211:PHE:HE2	1.71	0.56
1:M:225:PRO:HD2	1:M:228:TYR:HD2	1.71	0.56
1:O:253:ALA:O	1:O:256:VAL:HG23	2.05	0.56
2:H:270:GLU:OE2	2:H:274:ARG:NH1	2.38	0.56
1:E:108:THR:HG21	1:E:239:LEU:O	2.05	0.55
2:P:105:PHE:CE2	2:P:162:ARG:HG2	2.41	0.55
1:E:210:LEU:HB2	2:F:245:ASN:HB3	1.88	0.55
2:C:230:PHE:HB3	2:C:233:LEU:HD11	1.88	0.55
2:C:263:SER:OG	2:C:264:ALA:N	2.40	0.55
1:E:231:ILE:HG23	2:F:220:CYS:SG	2.46	0.55
1:G:46:GLY:HA2	1:G:52:MET:HE2	1.87	0.55
1:I:82:SER:HB3	1:I:85:LEU:H	1.71	0.55
1:I:137:VAL:HG22	2:J:156:THR:HG22	1.88	0.55
1:A:327:PHE:O	1:A:331:ILE:HG13	2.06	0.55
1:E:172:HIS:CD2	1:E:185:LEU:HD11	2.41	0.55
1:K:112:ASP:OD1	1:K:112:ASP:N	2.34	0.55
2:F:30:MET:HA	2:F:33:VAL:HG12	1.88	0.55
1:A:118:ILE:HD13	1:A:156:ALA:HA	1.89	0.55
2:J:119:THR:OG1	2:J:120:ARG:N	2.37	0.55
2:J:315:VAL:HG22	2:J:342:ILE:HD12	1.89	0.55
1:M:55:SER:O	1:M:58:LYS:N	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:116:GLY:HA3	2:N:320:LYS:HA	1.88	0.55
1:O:24:MET:HG2	1:O:36:TRP:CD1	2.41	0.55
1:B:189:ARG:NE	1:B:202:GLU:OE1	2.32	0.55
1:A:119:ARG:HD2	1:A:230:ASP:N	2.21	0.55
1:I:171:VAL:HB	1:I:224:MET:HE2	1.89	0.55
1:K:205:LEU:HD22	1:K:228:TYR:CG	2.42	0.55
1:K:324:CYS:O	1:K:328:THR:HG23	2.07	0.55
1:M:55:SER:O	1:M:59:GLU:HG2	2.06	0.55
2:L:116:GLY:HA3	2:L:320:LYS:HA	1.89	0.54
1:B:13:ASP:OD2	1:B:73:LYS:HB2	2.07	0.54
1:B:116:VAL:HB	1:B:221:VAL:HG22	1.89	0.54
2:L:325:ARG:HB2	2:L:331:GLY:HA3	1.90	0.54
2:P:107:ASN:HB2	2:P:132:GLN:OE1	2.07	0.54
2:C:343:CYS:O	2:C:347:LYS:HG3	2.07	0.54
2:F:81:ILE:HG21	2:F:95:ASP:HA	1.88	0.54
1:M:39:ARG:HE	1:M:56:GLU:HB3	1.71	0.54
1:G:204:TYR:O	1:G:208:VAL:HG23	2.08	0.54
1:M:307:THR:HG22	1:M:334:ARG:NH2	2.23	0.54
2:N:75:VAL:HG21	2:N:302:LEU:HD21	1.89	0.54
2:D:63:LEU:O	2:D:66:VAL:HG12	2.06	0.54
1:G:173:LYS:HB2	1:G:228:TYR:OH	2.08	0.54
2:P:182:VAL:HB	2:P:235:MET:HG2	1.90	0.54
2:D:18:THR:HG22	2:D:73:ASN:ND2	2.22	0.54
2:F:17:VAL:HG13	2:F:75:VAL:HG23	1.89	0.54
2:F:125:ASP:CG	2:F:176:ARG:HH22	2.15	0.54
1:G:157:PHE:O	1:G:161:ARG:HG3	2.06	0.54
1:O:242:GLY:HA2	2:P:221:MET:HE3	1.90	0.54
2:C:216:ILE:HD13	2:C:239:TYR:HB3	1.89	0.54
2:J:318:VAL:HG22	2:J:345:ARG:HD2	1.89	0.54
1:K:224:MET:HE2	1:K:229:GLY:N	2.23	0.54
2:L:236:PRO:HD2	2:L:239:TYR:HD2	1.73	0.54
1:O:186:GLN:O	1:O:190:GLU:HG3	2.08	0.54
1:B:54:PRO:HB3	1:G:56:GLU:OE1	2.08	0.54
1:B:324:CYS:O	1:B:328:THR:HG23	2.08	0.54
1:I:118:ILE:HD13	1:I:221:VAL:CG1	2.38	0.54
1:O:24:MET:HG2	1:O:36:TRP:NE1	2.23	0.54
1:O:273:ASP:OD1	1:O:325:SER:N	2.31	0.54
1:M:68:LEU:HD22	1:M:280:LEU:HD11	1.89	0.54
2:F:263:SER:HB3	2:F:266:TYR:HB2	1.90	0.54
1:E:232:LEU:HD12	1:E:235:LEU:HD23	1.90	0.53
1:E:332:CYS:O	1:E:336:LYS:HG3	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:109:PRO:O	2:J:120:ARG:HD3	2.07	0.53
1:O:157:PHE:CD2	1:O:191:VAL:HG12	2.43	0.53
1:G:171:VAL:HB	1:G:224:MET:HG2	1.91	0.53
2:H:260:GLU:HG3	2:H:262:TYR:CE2	2.43	0.53
1:O:239:LEU:HD23	2:P:224:VAL:HG11	1.88	0.53
2:D:16:PRO:HB2	2:D:73:ASN:OD1	2.08	0.53
1:E:273:ASP:OD2	1:E:325:SER:OG	2.25	0.53
1:I:206:ASP:OD1	1:I:206:ASP:N	2.42	0.53
1:K:120:GLU:OE1	1:K:122:THR:HB	2.08	0.53
1:M:301:GLU:HG2	1:M:305:PHE:CE2	2.44	0.53
2:L:236:PRO:HD2	2:L:239:TYR:CD2	2.44	0.53
2:L:263:SER:OG	2:L:264:ALA:N	2.37	0.53
2:N:176:ARG:NH1	2:N:231:ASP:OD1	2.40	0.53
1:B:9:LEU:HD23	1:B:11:PRO:HG3	1.90	0.53
1:I:261:SER:HB2	1:I:263:HIS:CD2	2.44	0.53
1:O:109:PRO:HB3	2:P:120:ARG:HB3	1.90	0.53
2:C:325:ARG:HB2	2:C:331:GLY:HA3	1.90	0.53
1:G:155:PHE:CZ	1:G:251:ILE:HG13	2.43	0.53
1:O:314:LEU:HB3	1:O:320:GLY:HA3	1.90	0.53
2:F:116:GLY:O	2:F:325:ARG:NH2	2.37	0.53
2:H:75:VAL:HG21	2:H:302:LEU:HD21	1.91	0.53
1:I:108:THR:HG21	1:I:239:LEU:O	2.09	0.53
1:K:206:ASP:OD1	1:K:206:ASP:N	2.40	0.53
1:G:235:LEU:CD1	2:H:224:VAL:HG21	2.38	0.52
1:M:66:MET:HG3	1:M:291:MET:HE3	1.91	0.52
2:P:263:SER:OG	2:P:264:ALA:N	2.41	0.52
1:K:165:ARG:HD2	1:K:220:ASP:OD1	2.09	0.52
1:M:165:ARG:HD2	1:M:220:ASP:OD1	2.09	0.52
1:O:169:THR:HG21	1:O:203:MET:HE3	1.92	0.52
1:I:313:SER:O	1:I:327:PHE:HD1	1.92	0.52
1:K:227:LEU:O	1:K:231:ILE:HG12	2.08	0.52
1:G:243:LEU:HB3	1:G:262:VAL:HG23	1.92	0.52
2:H:310:MET:HG2	2:H:346:VAL:HG23	1.90	0.52
1:I:231:ILE:HD12	2:J:216:ILE:HG13	1.92	0.52
2:F:105:PHE:CE2	2:F:162:ARG:HG2	2.45	0.52
1:G:88:ARG:HD2	1:G:121:ASN:O	2.09	0.52
1:G:109:PRO:HB3	2:H:120:ARG:HG2	1.90	0.52
1:I:126:TYR:OH	2:J:184:LYS:HD3	2.09	0.52
1:O:24:MET:HE3	1:O:36:TRP:CD1	2.45	0.52
1:B:214:GLN:O	2:C:120:ARG:NH2	2.43	0.52
1:A:296:HIS:O	1:A:300:ILE:HG13	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:287:ASN:HB2	2:D:328:ASP:OD1	2.10	0.52
2:N:176:ARG:HD3	2:N:231:ASP:OD1	2.09	0.52
2:H:134:GLU:HG3	2:H:159:LYS:HD2	1.90	0.52
2:H:180:THR:HB	2:H:233:LEU:HD23	1.91	0.52
1:I:22:ALA:O	1:I:26:ILE:HG13	2.10	0.52
1:A:276:ASN:HD22	1:A:318:LEU:HD13	1.73	0.52
1:E:69:LYS:HE2	1:E:260:GLU:HB3	1.91	0.52
1:O:157:PHE:CE2	1:O:192:ALA:HA	2.44	0.52
2:C:142:HIS:NE2	2:D:150:GLU:OE2	2.33	0.52
2:H:236:PRO:HD2	2:H:239:TYR:HD2	1.74	0.52
1:O:97:VAL:HG13	1:O:118:ILE:HG22	1.92	0.52
1:I:43:ALA:HB3	1:I:80:HIS:CE1	2.45	0.52
2:J:119:THR:HG23	2:J:121:HIS:H	1.74	0.52
1:K:63:LYS:HD3	1:K:63:LYS:C	2.34	0.52
2:D:18:THR:CG2	2:D:73:ASN:HD22	2.22	0.51
2:J:18:THR:CG2	2:J:69:SER:HB3	2.40	0.51
2:P:310:MET:HG2	2:P:346:VAL:HG23	1.91	0.51
1:M:58:LYS:HE2	1:M:62:ASP:OD2	2.10	0.51
1:M:14:GLY:HA3	1:M:265:THR:HB	1.92	0.51
1:A:206:ASP:OD1	1:A:206:ASP:N	2.43	0.51
2:H:334:THR:O	2:H:337:ASP:HB2	2.10	0.51
1:M:235:LEU:HD13	2:N:246:LEU:HD11	1.93	0.51
2:P:221:MET:HE2	2:P:225:GLN:NE2	2.26	0.51
1:I:108:THR:HG22	1:I:110:TYR:H	1.74	0.51
2:L:20:LEU:HB2	2:L:78:ILE:HG22	1.92	0.51
1:A:268:ASP:N	1:A:268:ASP:OD1	2.43	0.51
1:M:132:VAL:HG22	1:M:138:VAL:HG22	1.93	0.51
1:M:304:CYS:O	1:M:308:ILE:HG13	2.11	0.51
1:O:225:PRO:HD2	1:O:228:TYR:HD2	1.76	0.51
1:A:160:ALA:HA	1:A:165:ARG:HB2	1.92	0.51
1:A:332:CYS:O	1:A:336:LYS:HD2	2.10	0.51
1:I:146:GLU:HB2	1:I:187:LYS:HE3	1.92	0.51
2:J:18:THR:HG21	2:J:69:SER:HB3	1.93	0.51
1:E:273:ASP:OD2	1:E:323:LYS:HE2	2.11	0.51
1:B:273:ASP:OD2	1:B:325:SER:OG	2.23	0.51
1:A:330:GLU:HG3	1:A:333:ARG:NH1	2.25	0.51
2:F:263:SER:OG	2:F:264:ALA:N	2.44	0.51
2:J:169:ASP:O	2:J:173:LYS:HB2	2.11	0.51
2:L:19:MET:HE1	2:L:34:LYS:HD3	1.93	0.51
1:M:244:GLY:O	1:M:279:ALA:HB2	2.11	0.51
2:N:40:ALA:HB1	2:N:310:MET:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:136:GLU:OE2	2:N:136:GLU:N	2.34	0.51
2:P:172:THR:HG22	2:P:209:ILE:HD11	1.92	0.51
2:C:275:HIS:NE2	2:C:278:ALA:HA	2.26	0.51
2:P:131:GLU:OE1	2:P:133:THR:OG1	2.27	0.51
2:P:247:ALA:O	2:P:250:LEU:HB2	2.11	0.50
2:P:342:ILE:O	2:P:346:VAL:HG12	2.11	0.50
2:P:40:ALA:HB1	2:P:310:MET:HE3	1.92	0.50
1:K:210:LEU:HB2	2:L:245:ASN:HB3	1.93	0.50
2:L:110:HIS:NE2	2:L:260:GLU:OE1	2.44	0.50
2:L:341:GLU:OE1	2:L:341:GLU:C	2.54	0.50
2:C:260:GLU:HG2	2:C:269:PHE:CD2	2.46	0.50
1:A:276:ASN:ND2	1:A:318:LEU:CD1	2.72	0.50
2:F:142:HIS:NE2	2:H:150:GLU:OE2	2.32	0.50
2:H:70:MET:SD	2:H:76:ALA:HB2	2.52	0.50
1:M:6:THR:HG22	1:M:35:GLN:HG3	1.92	0.50
1:O:69:LYS:HB3	1:O:260:GLU:HB3	1.94	0.50
1:O:172:HIS:CE1	1:O:185:LEU:HD11	2.47	0.50
1:A:225:PRO:HD2	1:A:228:TYR:CD2	2.46	0.50
2:J:14:SER:OG	2:J:15:PHE:N	2.30	0.50
1:M:118:ILE:HD13	1:M:221:VAL:HG13	1.93	0.50
1:M:137:VAL:HG22	2:N:156:THR:HG22	1.94	0.50
1:O:138:VAL:HG11	2:P:190:LEU:HB3	1.93	0.50
2:P:115:PRO:HB2	2:P:320:LYS:HD2	1.93	0.50
1:M:119:ARG:HD2	1:M:230:ASP:N	2.27	0.50
1:O:14:GLY:HA3	1:O:265:THR:HG23	1.94	0.50
1:E:63:LYS:O	1:E:65:LYS:N	2.39	0.50
1:K:268:ASP:OD1	1:K:268:ASP:N	2.45	0.50
2:C:104:LEU:HD11	2:C:268:VAL:HG21	1.94	0.49
1:A:292:GLY:HA2	1:A:294:PHE:CE1	2.46	0.49
1:E:177:MET:HG2	2:F:151:CYS:SG	2.52	0.49
1:G:206:ASP:N	1:G:206:ASP:OD1	2.46	0.49
2:H:342:ILE:O	2:H:346:VAL:HG12	2.12	0.49
1:M:132:VAL:HG23	2:N:190:LEU:HD13	1.93	0.49
1:M:206:ASP:OD2	2:N:275:HIS:HA	2.12	0.49
1:M:330:GLU:CD	1:M:334:ARG:HH22	2.20	0.49
2:H:328:ASP:OD1	2:H:328:ASP:N	2.44	0.49
1:A:114:ASN:OD1	1:A:165:ARG:NH2	2.35	0.49
1:E:142:LYS:NZ	1:E:181:ASP:OD1	2.43	0.49
1:I:330:GLU:OE2	1:I:334:ARG:NH1	2.46	0.49
2:N:200:GLU:HG3	2:N:211:PHE:CE2	2.47	0.49
1:O:64:ASN:C	1:O:66:MET:H	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:8:THR:OG1	1:M:60:SER:OG	2.23	0.49
1:O:189:ARG:NE	1:O:202:GLU:OE1	2.46	0.49
1:O:332:CYS:HA	1:O:335:VAL:HG12	1.94	0.49
2:P:263:SER:HB3	2:P:266:TYR:H	1.77	0.49
2:H:111:VAL:HG21	2:H:128:ILE:HD11	1.95	0.49
1:K:114:ASN:OD1	1:K:165:ARG:NH2	2.44	0.49
2:N:40:ALA:HB3	2:N:42:VAL:HG23	1.94	0.49
2:C:40:ALA:HB1	2:C:310:MET:HE3	1.95	0.49
2:C:134:GLU:OE1	2:C:156:THR:OG1	2.29	0.49
1:A:125:GLU:HG2	1:A:142:LYS:HB2	1.95	0.49
2:F:40:ALA:HB3	2:F:42:VAL:HG23	1.94	0.49
1:I:58:LYS:NZ	1:I:62:ASP:OD2	2.46	0.49
1:M:26:ILE:HD11	1:M:328:THR:HG23	1.95	0.49
2:N:226:ASN:O	2:N:229:GLN:HG2	2.12	0.49
1:O:108:THR:C	1:O:110:TYR:H	2.21	0.49
1:B:130:GLU:OE1	1:B:140:SER:OG	2.30	0.49
1:I:126:TYR:OH	1:I:227:LEU:HD12	2.13	0.49
2:P:260:GLU:HG2	2:P:269:PHE:CE2	2.48	0.49
1:G:88:ARG:HH11	1:G:88:ARG:HG2	1.78	0.48
2:J:116:GLY:O	2:J:325:ARG:NH2	2.46	0.48
1:O:6:THR:O	1:O:64:ASN:HB3	2.12	0.48
2:J:67:LEU:HD13	2:J:102:LEU:HD23	1.95	0.48
1:G:276:ASN:HD21	1:G:318:LEU:HD11	1.78	0.48
1:I:204:TYR:CD2	2:J:276:PRO:HG2	2.48	0.48
2:J:275:HIS:O	2:J:279:GLN:NE2	2.46	0.48
1:K:224:MET:CE	1:K:228:TYR:C	2.87	0.48
2:N:263:SER:HB3	2:N:266:TYR:HB2	1.93	0.48
1:O:16:GLY:O	1:O:20:SER:OG	2.13	0.48
1:B:155:PHE:CZ	1:B:251:ILE:HG13	2.48	0.48
1:A:272:LYS:HG2	1:A:274:MET:HE3	1.94	0.48
2:F:20:LEU:HB2	2:F:78:ILE:HG22	1.95	0.48
2:J:325:ARG:HB2	2:J:331:GLY:HA3	1.94	0.48
2:N:197:GLN:O	2:N:201:GLU:HG3	2.13	0.48
1:O:27:PHE:HD2	1:O:34:ILE:CD1	2.27	0.48
2:H:263:SER:OG	2:H:264:ALA:N	2.46	0.48
2:N:296:SER:HB3	2:N:311:ILE:HD11	1.94	0.48
1:A:330:GLU:O	1:A:334:ARG:HG3	2.14	0.48
1:G:179:MET:H	2:H:141:GLU:CD	2.22	0.48
1:I:119:ARG:HD2	1:I:229:GLY:C	2.39	0.48
1:M:225:PRO:HD2	1:M:228:TYR:CD2	2.49	0.48
1:A:250:ASN:ND2	1:A:260:GLU:OE2	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:88:ARG:HH11	1:E:88:ARG:HG2	1.79	0.48
1:M:275:ALA:O	1:M:324:CYS:HB2	2.14	0.48
2:N:107:ASN:HB2	2:N:132:GLN:CD	2.39	0.48
1:E:189:ARG:NE	1:E:202:GLU:OE1	2.43	0.48
2:J:42:VAL:HG12	2:J:44:VAL:HG23	1.95	0.48
1:B:37:GLU:CG	1:B:39:ARG:NH1	2.75	0.48
2:L:18:THR:HG23	2:L:76:ALA:HB2	1.95	0.48
1:O:123:GLU:HB2	1:O:143:LEU:O	2.14	0.48
1:O:126:TYR:OH	2:P:184:LYS:HE2	2.14	0.48
1:G:137:VAL:HG22	2:H:156:THR:HG22	1.96	0.48
2:H:208:LYS:HD2	2:H:208:LYS:HA	1.59	0.48
2:P:221:MET:HE2	2:P:225:GLN:HE22	1.77	0.48
1:B:307:THR:HA	1:B:334:ARG:NH1	2.27	0.47
1:A:69:LYS:O	1:A:260:GLU:HA	2.14	0.47
1:A:315:THR:OG1	1:A:317:ASP:OD1	2.25	0.47
1:I:30:ALA:HB2	1:I:335:VAL:HG21	1.96	0.47
2:J:19:MET:HB3	2:J:48:GLU:HG3	1.95	0.47
2:N:285:ILE:HG13	2:N:285:ILE:O	2.14	0.47
1:O:119:ARG:HD2	1:O:230:ASP:N	2.29	0.47
2:H:165:LYS:HA	2:H:202:VAL:HG11	1.95	0.47
2:J:18:THR:HG22	2:J:73:ASN:ND2	2.21	0.47
1:K:94:TYR:CE2	1:K:151:ARG:HG2	2.49	0.47
1:K:231:ILE:HD12	2:L:216:ILE:HG13	1.96	0.47
1:M:143:LEU:HD21	2:P:144:SER:O	2.14	0.47
1:O:9:LEU:O	1:O:11:PRO:HD3	2.15	0.47
2:C:156:THR:OG1	2:C:159:LYS:HG3	2.14	0.47
2:H:19:MET:O	2:H:21:PRO:HD3	2.13	0.47
1:I:126:TYR:CD1	2:J:187:ILE:HG21	2.48	0.47
1:I:224:MET:HE2	1:I:224:MET:HB3	1.71	0.47
1:B:149:SER:O	1:B:152:ILE:HG22	2.15	0.47
2:F:314:ALA:CB	2:F:346:VAL:HG23	2.44	0.47
2:H:287:ASN:HD21	2:H:329:MET:HE2	1.80	0.47
2:J:71:LYS:O	2:J:74:LYS:NZ	2.47	0.47
1:A:130:GLU:OE1	1:A:140:SER:OG	2.33	0.47
1:E:119:ARG:HD2	1:E:230:ASP:N	2.28	0.47
1:I:165:ARG:NH1	1:I:220:ASP:OD1	2.38	0.47
1:I:288:LEU:HB2	1:I:297:ALA:HB2	1.97	0.47
1:K:98:ARG:HD2	1:K:233:SER:HB2	1.97	0.47
1:O:117:THR:HB	1:O:224:MET:HE1	1.96	0.47
2:P:103:ASP:CG	2:P:162:ARG:HH22	2.22	0.47
1:A:208:VAL:O	1:A:212:MET:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:CYS:O	1:A:335:VAL:HG12	2.15	0.47
1:E:97:VAL:HG13	1:E:118:ILE:HG13	1.96	0.47
1:K:206:ASP:HB3	2:L:241:ASN:OD1	2.14	0.47
2:N:25:VAL:HG21	2:N:286:ALA:CB	2.44	0.47
2:P:153:LYS:NZ	2:P:192:ASP:OD1	2.36	0.47
2:C:338:PHE:O	2:C:342:ILE:HG12	2.15	0.47
2:D:317:LYS:HE2	2:D:345:ARG:HD2	1.96	0.47
1:E:215:ASP:O	1:E:218:GLN:HG2	2.15	0.47
2:F:315:VAL:HG22	2:F:342:ILE:HD13	1.96	0.47
1:G:215:ASP:OD1	1:G:217:SER:HB3	2.14	0.47
2:L:345:ARG:HA	2:L:345:ARG:NE	2.30	0.47
2:N:34:LYS:NZ	2:N:48:GLU:OE2	2.33	0.47
2:P:16:PRO:HB2	2:P:73:ASN:ND2	2.30	0.47
1:B:334:ARG:HG3	1:B:334:ARG:NH1	2.28	0.47
2:F:125:ASP:OD1	2:F:176:ARG:NH2	2.42	0.47
2:J:118:MET:HE3	2:J:122:ASN:ND2	2.30	0.47
1:K:178:ARG:HH12	2:L:93:SER:HB3	1.80	0.47
2:N:311:ILE:O	2:N:315:VAL:HG23	2.15	0.47
1:O:291:MET:HB2	1:O:293:LEU:HG	1.97	0.47
2:D:172:THR:HG22	2:D:209:ILE:HD11	1.97	0.46
1:G:93:LEU:HD23	1:G:252:GLY:HA3	1.98	0.46
1:G:292:GLY:HA2	1:G:294:PHE:HE1	1.80	0.46
1:I:104:GLU:CD	1:I:309:LYS:HD2	2.40	0.46
1:I:205:LEU:HD13	1:I:228:TYR:HB3	1.96	0.46
1:K:179:MET:HE3	1:K:179:MET:HB2	1.79	0.46
1:M:313:SER:O	1:M:313:SER:OG	2.21	0.46
2:P:105:PHE:CZ	2:P:162:ARG:HG2	2.51	0.46
2:C:155:VAL:HG12	2:C:194:LEU:HD23	1.96	0.46
2:F:70:MET:SD	2:F:76:ALA:HB2	2.55	0.46
1:M:123:GLU:HB2	1:M:143:LEU:O	2.15	0.46
2:P:235:MET:HB2	2:P:235:MET:HE3	1.32	0.46
2:C:210:LYS:HE3	2:C:212:GLU:HG3	1.98	0.46
1:G:150:LYS:O	1:G:154:GLU:HG3	2.15	0.46
1:I:225:PRO:HD2	1:I:228:TYR:CD2	2.50	0.46
1:O:27:PHE:CD2	1:O:34:ILE:HD11	2.49	0.46
2:C:115:PRO:HB2	2:C:320:LYS:HD2	1.96	0.46
2:H:113:SER:HB2	2:H:124:LEU:H	1.80	0.46
2:J:96:MET:HE1	2:J:133:THR:O	2.15	0.46
2:J:128:ILE:HD13	2:J:233:LEU:HD13	1.98	0.46
1:B:22:ALA:O	1:B:26:ILE:HG13	2.15	0.46
2:D:263:SER:OG	2:D:264:ALA:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:296:SER:O	2:F:308:SER:OG	2.26	0.46
1:M:113:VAL:HG22	1:M:216:PRO:HB2	1.98	0.46
2:P:66:VAL:O	2:P:69:SER:HB2	2.15	0.46
2:C:14:SER:OG	2:C:15:PHE:N	2.47	0.46
2:C:93:SER:OG	2:C:94:TYR:N	2.48	0.46
1:E:204:TYR:HD2	2:F:277:PHE:HE2	1.64	0.46
2:H:236:PRO:HD2	2:H:239:TYR:CD2	2.50	0.46
2:J:183:HIS:O	2:J:215:ILE:HA	2.16	0.46
1:M:15:ILE:HG12	1:M:265:THR:O	2.16	0.46
1:O:126:TYR:CG	2:P:187:ILE:HG21	2.50	0.46
1:O:287:MET:HG2	1:O:291:MET:HE2	1.97	0.46
2:C:76:ALA:HB3	2:C:268:VAL:HG13	1.98	0.46
2:F:157:ARG:NH2	2:F:161:GLN:NE2	2.64	0.46
2:P:317:LYS:HD2	2:P:317:LYS:N	2.29	0.46
1:B:119:ARG:HD2	1:B:229:GLY:C	2.41	0.46
1:G:206:ASP:HB3	2:H:241:ASN:ND2	2.30	0.46
2:J:269:PHE:CZ	2:J:298:MET:HA	2.51	0.46
1:K:19:ILE:HG13	1:K:280:LEU:HD22	1.98	0.46
1:M:112:ASP:OD1	1:M:112:ASP:N	2.49	0.46
1:B:131:HIS:HB3	2:D:140:LEU:HD22	1.97	0.46
1:G:22:ALA:O	1:G:26:ILE:HG13	2.16	0.46
2:N:65:GLN:OE1	2:N:65:GLN:N	2.38	0.46
2:C:180:THR:HA	2:C:212:GLU:O	2.16	0.45
1:A:184:PHE:CE2	1:A:225:PRO:HD3	2.51	0.45
1:M:86:LEU:O	1:M:90:THR:HG22	2.16	0.45
2:N:125:ASP:CG	2:N:176:ARG:HH22	2.24	0.45
1:O:311:GLY:HA2	1:O:314:LEU:HD22	1.98	0.45
1:B:165:ARG:NH1	1:B:220:ASP:OD1	2.44	0.45
2:H:132:GLN:HE22	2:H:261:SER:HB3	1.80	0.45
2:J:15:PHE:HB2	2:J:43:PRO:O	2.16	0.45
1:K:171:VAL:HB	1:K:224:MET:CG	2.47	0.45
2:L:182:VAL:HB	2:L:235:MET:CG	2.45	0.45
1:B:129:ILE:CG2	1:B:141:ILE:HG13	2.47	0.45
1:I:222:LEU:HB3	1:I:224:MET:HE1	1.97	0.45
1:M:329:GLU:HA	1:M:332:CYS:HB2	1.98	0.45
1:O:109:PRO:O	2:P:120:ARG:HD3	2.16	0.45
1:G:288:LEU:HB3	1:G:297:ALA:HB2	1.97	0.45
1:I:143:LEU:HD21	2:L:144:SER:O	2.16	0.45
1:M:186:GLN:O	1:M:190:GLU:HG3	2.15	0.45
1:O:288:LEU:HA	1:O:291:MET:HG3	1.98	0.45
1:A:80:HIS:HE1	1:A:83:MET:HE2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:232:LEU:HD12	1:M:232:LEU:HA	1.75	0.45
2:C:19:MET:HG3	2:C:77:ILE:HG23	1.99	0.45
1:E:179:MET:HG2	2:F:149:ILE:CG2	2.46	0.45
2:F:136:GLU:OE2	2:F:136:GLU:N	2.42	0.45
1:I:143:LEU:HD12	1:I:143:LEU:HA	1.82	0.45
1:I:222:LEU:HB3	1:I:224:MET:CE	2.45	0.45
2:L:49:HIS:NE2	2:L:69:SER:HB2	2.32	0.45
2:L:116:GLY:C	2:L:325:ARG:HH22	2.25	0.45
1:B:126:TYR:CD1	2:C:187:ILE:HG21	2.51	0.45
2:L:216:ILE:HD13	2:L:239:TYR:HB3	1.99	0.45
2:P:15:PHE:CE1	2:P:304:LEU:HD11	2.51	0.45
1:B:30:ALA:O	1:B:299:ARG:NH2	2.50	0.45
1:B:129:ILE:HG22	1:B:141:ILE:CG1	2.47	0.45
1:G:224:MET:HE3	1:G:224:MET:HB2	1.85	0.45
1:E:137:VAL:HG22	2:F:156:THR:HG22	1.98	0.45
1:E:206:ASP:OD1	1:E:206:ASP:N	2.49	0.45
1:K:105:GLY:HA3	1:K:308:ILE:HG22	1.98	0.45
1:M:6:THR:HG21	1:M:35:GLN:NE2	2.32	0.45
1:O:259:PHE:CZ	1:O:287:MET:HA	2.52	0.45
1:B:165:ARG:HD2	1:B:220:ASP:OD1	2.17	0.45
1:A:161:ARG:HH12	1:A:197:ASP:CG	2.25	0.45
2:F:144:SER:HB3	2:H:150:GLU:OE2	2.17	0.45
2:H:71:LYS:O	2:H:74:LYS:HE2	2.17	0.45
1:I:24:MET:HG2	1:I:36:TRP:NE1	2.32	0.45
1:I:130:GLU:HB3	1:I:140:SER:HA	1.97	0.45
2:J:157:ARG:O	2:J:161:GLN:HG2	2.17	0.45
1:M:205:LEU:HD22	1:M:228:TYR:CD1	2.52	0.45
2:C:294:SER:HA	2:C:297:ASN:HB2	1.99	0.44
1:A:109:PRO:O	2:D:120:ARG:HD3	2.16	0.44
1:E:232:LEU:HD12	1:E:232:LEU:HA	1.82	0.44
1:I:15:ILE:HD13	1:I:266:ALA:HB3	1.99	0.44
2:L:107:ASN:HB2	2:L:132:GLN:CD	2.42	0.44
2:L:179:VAL:O	2:L:211:PHE:HA	2.17	0.44
1:M:307:THR:HG22	1:M:334:ARG:HH21	1.81	0.44
1:A:112:ASP:OD1	1:A:112:ASP:N	2.45	0.44
2:D:108:VAL:HG22	2:D:129:ILE:HG12	1.98	0.44
2:F:336:HIS:NE2	2:J:336:HIS:CE1	2.86	0.44
1:K:130:GLU:OE1	1:K:140:SER:OG	2.33	0.44
2:N:336:HIS:NE2	2:N:340:GLU:OE2	2.50	0.44
2:C:210:LYS:HG2	2:C:211:PHE:N	2.33	0.44
1:E:160:ALA:HA	1:E:165:ARG:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:130:ARG:HG3	2:H:240:GLY:HA3	1.98	0.44
1:M:23:VAL:HG23	1:M:281:LEU:CD1	2.47	0.44
2:D:14:SER:OG	2:D:15:PHE:N	2.50	0.44
1:G:99:PRO:HA	1:G:116:VAL:HA	1.99	0.44
1:M:303:ALA:O	1:M:307:THR:HG23	2.18	0.44
2:C:70:MET:HE3	2:C:70:MET:HB3	1.82	0.44
2:F:63:LEU:HG	2:F:65:GLN:CD	2.42	0.44
2:H:78:ILE:HG12	2:H:268:VAL:HG12	1.98	0.44
2:J:318:VAL:HG21	2:J:342:ILE:HD13	1.99	0.44
1:K:38:GLU:HB2	1:M:45:GLN:HG3	1.98	0.44
2:N:153:LYS:NZ	2:N:192:ASP:OD1	2.43	0.44
1:O:123:GLU:HB2	1:O:124:GLY:H	1.62	0.44
2:D:99:ARG:NH2	2:D:270:GLU:OE1	2.51	0.44
2:D:126:LEU:HD22	2:D:230:PHE:HB2	1.98	0.44
2:D:153:LYS:NZ	2:D:192:ASP:OD1	2.41	0.44
2:D:275:HIS:NE2	2:D:278:ALA:HA	2.33	0.44
1:G:232:LEU:HD12	1:G:232:LEU:HA	1.83	0.44
1:O:281:LEU:HD23	1:O:281:LEU:HA	1.71	0.44
1:A:137:VAL:HG22	2:D:156:THR:HG22	2.00	0.44
2:H:129:ILE:HG22	2:H:163:ILE:HD11	2.00	0.44
2:L:189:LYS:O	2:L:193:GLY:HA3	2.17	0.44
2:C:292:LEU:HD23	2:C:292:LEU:HA	1.84	0.44
1:E:142:LYS:HB3	2:F:151:CYS:HB3	2.00	0.44
1:I:332:CYS:HA	1:I:335:VAL:HG12	1.99	0.44
1:K:123:GLU:HB2	1:K:124:GLY:H	1.58	0.44
2:L:67:LEU:HD23	2:L:67:LEU:O	2.18	0.44
1:M:69:LYS:O	1:M:260:GLU:HA	2.18	0.44
2:N:263:SER:OG	2:N:264:ALA:N	2.50	0.44
2:D:34:LYS:HE2	2:D:34:LYS:HB2	1.78	0.44
1:O:284:ALA:O	1:O:287:MET:HB3	2.18	0.44
1:B:141:ILE:HD13	2:C:152:LEU:HD22	1.99	0.43
1:B:206:ASP:N	1:B:206:ASP:OD1	2.48	0.43
1:G:244:GLY:O	1:G:279:ALA:HB2	2.18	0.43
1:K:21:ALA:HA	1:K:24:MET:HE3	2.00	0.43
2:L:32:ALA:O	2:L:36:VAL:HG23	2.17	0.43
2:L:300:ARG:O	2:L:300:ARG:HG2	2.18	0.43
1:M:65:LYS:C	1:M:66:MET:HE2	2.43	0.43
1:M:119:ARG:HD2	1:M:229:GLY:C	2.42	0.43
1:O:125:GLU:C	1:O:126:TYR:HD1	2.26	0.43
2:P:99:ARG:HD2	2:P:132:GLN:NE2	2.33	0.43
2:C:293:LEU:HA	2:C:293:LEU:HD23	1.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:102:LEU:N	2:D:102:LEU:HD12	2.33	0.43
2:H:171:ALA:HA	2:H:176:ARG:HB2	2.00	0.43
1:K:126:TYR:CD1	2:L:187:ILE:HG21	2.53	0.43
1:B:6:THR:HG22	1:B:35:GLN:HB2	2.00	0.43
1:A:157:PHE:O	1:A:161:ARG:HG3	2.18	0.43
2:F:157:ARG:NH2	2:F:161:GLN:HE22	2.17	0.43
2:J:134:GLU:OE2	2:J:156:THR:N	2.51	0.43
2:L:230:PHE:CD1	2:L:233:LEU:HD21	2.53	0.43
1:M:39:ARG:NE	1:M:56:GLU:O	2.51	0.43
1:M:210:LEU:HB2	2:N:245:ASN:HB3	1.99	0.43
1:O:119:ARG:HG3	1:O:229:GLY:HA3	1.99	0.43
1:O:276:ASN:HA	1:O:277:PRO:HD2	1.80	0.43
1:B:277:PRO:HG3	1:B:328:THR:HG22	2.01	0.43
1:E:86:LEU:HD12	1:E:86:LEU:HA	1.80	0.43
1:G:161:ARG:NH1	1:G:197:ASP:OD1	2.51	0.43
1:K:285:VAL:HG13	1:K:297:ALA:HB1	1.99	0.43
2:N:81:ILE:HG21	2:N:95:ASP:CB	2.46	0.43
1:O:145:THR:HG22	2:P:148:VAL:HG13	2.00	0.43
1:O:289:ARG:HG2	1:O:289:ARG:HH11	1.83	0.43
1:B:224:MET:HE2	1:B:224:MET:HB3	1.72	0.43
1:A:43:ALA:H	1:A:83:MET:HE1	1.83	0.43
2:D:133:THR:OG1	2:D:159:LYS:HD3	2.18	0.43
1:E:160:ALA:HB3	1:E:198:ILE:HD13	1.99	0.43
1:G:171:VAL:HB	1:G:224:MET:CG	2.47	0.43
1:M:145:THR:HG22	2:N:148:VAL:HG13	2.01	0.43
1:O:232:LEU:HD12	1:O:232:LEU:HA	1.83	0.43
2:P:165:LYS:HA	2:P:202:VAL:HG11	2.00	0.43
1:B:33:PRO:HG2	1:B:293:LEU:HD13	1.99	0.43
2:D:30:MET:HE1	2:D:77:ILE:HG12	2.01	0.43
1:G:288:LEU:CB	1:G:297:ALA:HB2	2.48	0.43
1:I:246:THR:HA	1:I:247:PRO:HD3	1.91	0.43
2:J:180:THR:O	2:J:233:LEU:HA	2.17	0.43
1:K:307:THR:HG23	1:K:330:GLU:OE2	2.18	0.43
1:M:231:ILE:HD13	2:N:242:ILE:CD1	2.47	0.43
1:B:251:ILE:HD12	1:B:290:HIS:NE2	2.34	0.43
1:E:98:ARG:HD2	1:E:233:SER:HB2	2.00	0.43
2:N:205:LEU:C	2:N:207:PRO:HD3	2.44	0.43
1:O:86:LEU:HD23	1:O:86:LEU:HA	1.80	0.43
1:O:259:PHE:CD1	1:O:287:MET:HB2	2.54	0.43
2:C:107:ASN:HB2	2:C:132:GLN:CD	2.44	0.43
2:D:47:GLN:HE22	2:D:73:ASN:HD21	1.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:70:MET:CE	2:F:75:VAL:C	2.92	0.43
1:G:292:GLY:HA2	1:G:294:PHE:CE1	2.54	0.43
1:G:332:CYS:O	1:G:335:VAL:HG12	2.19	0.43
2:L:296:SER:OG	2:L:312:ALA:HB2	2.19	0.43
1:A:161:ARG:NH1	1:A:197:ASP:CG	2.76	0.43
1:G:287:MET:HG2	1:G:291:MET:HE2	2.00	0.43
1:I:317:ASP:OD2	1:I:318:LEU:HD13	2.19	0.43
2:L:110:HIS:NE2	2:L:260:GLU:CD	2.77	0.43
1:M:41:VAL:HG22	1:M:53:ILE:HD12	1.99	0.43
1:O:283:SER:O	1:O:286:MET:HB2	2.18	0.43
1:B:295:ASP:O	1:B:299:ARG:HG3	2.19	0.43
1:G:285:VAL:HG13	1:G:297:ALA:HB1	1.99	0.43
2:J:216:ILE:HD13	2:J:239:TYR:HB3	1.99	0.43
1:O:284:ALA:HA	1:O:287:MET:HE2	2.01	0.43
1:O:327:PHE:CE2	1:O:331:ILE:HD11	2.54	0.43
2:P:95:ASP:O	2:P:98:LEU:HB3	2.19	0.43
1:B:6:THR:O	1:B:64:ASN:HB3	2.19	0.42
1:B:16:GLY:HA2	1:B:71:PRO:HD2	2.01	0.42
1:B:126:TYR:CE1	1:B:227:LEU:HD12	2.54	0.42
1:B:181:ASP:OD2	1:B:228:TYR:OH	2.36	0.42
1:E:161:ARG:HD3	1:E:197:ASP:OD2	2.19	0.42
1:G:204:TYR:CD2	2:H:276:PRO:HG2	2.54	0.42
1:O:328:THR:HA	1:O:331:ILE:HD12	2.01	0.42
2:C:315:VAL:HG22	2:C:342:ILE:HD12	2.00	0.42
2:D:223:LEU:HD12	2:D:230:PHE:CE2	2.54	0.42
1:I:5:GLN:O	1:I:34:ILE:HA	2.19	0.42
2:L:324:VAL:CG1	2:L:338:PHE:HA	2.45	0.42
2:N:235:MET:HE2	2:N:235:MET:HB3	1.80	0.42
1:O:204:TYR:CD2	2:P:276:PRO:HG2	2.54	0.42
1:O:314:LEU:HD12	1:O:314:LEU:HA	1.65	0.42
1:A:119:ARG:HD2	1:A:229:GLY:C	2.45	0.42
1:E:88:ARG:HG2	1:E:88:ARG:NH1	2.34	0.42
1:G:61:MET:CE	1:G:67:GLY:HA3	2.50	0.42
1:G:243:LEU:O	1:G:263:HIS:HB3	2.18	0.42
1:G:292:GLY:O	1:G:294:PHE:HD1	2.02	0.42
1:K:280:LEU:HD12	1:K:280:LEU:HA	1.86	0.42
2:L:129:ILE:O	2:L:234:VAL:HA	2.19	0.42
1:M:43:ALA:HB2	1:M:83:MET:HE3	2.01	0.42
1:M:236:CYS:HA	1:M:239:LEU:HD12	2.02	0.42
1:O:119:ARG:HG2	1:O:120:GLU:O	2.19	0.42
1:A:165:ARG:HD2	1:A:220:ASP:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:70:MET:HE1	2:F:75:VAL:C	2.45	0.42
1:I:248:SER:N	1:I:286:MET:HE1	2.35	0.42
2:C:171:ALA:HA	2:C:176:ARG:HB2	2.01	0.42
1:A:113:VAL:HG22	1:A:216:PRO:HB2	2.00	0.42
1:E:204:TYR:CD2	2:F:276:PRO:HG2	2.54	0.42
2:J:317:LYS:HG2	2:J:345:ARG:NH1	2.35	0.42
2:N:42:VAL:HG22	2:N:307:HIS:HB3	2.01	0.42
1:O:225:PRO:HD2	1:O:228:TYR:CD2	2.53	0.42
2:P:196:LEU:HD23	2:P:196:LEU:HA	1.70	0.42
1:B:37:GLU:CD	1:B:39:ARG:NH1	2.77	0.42
1:B:161:ARG:NH1	1:B:197:ASP:OD1	2.52	0.42
1:A:210:LEU:HB2	2:D:245:ASN:HB3	2.02	0.42
1:E:47:PRO:HD3	1:E:52:MET:CE	2.49	0.42
2:L:160:SER:HA	2:L:163:ILE:HG22	2.00	0.42
2:L:300:ARG:NH1	2:L:305:GLU:OE2	2.52	0.42
1:O:77:ALA:HB3	1:O:80:HIS:HB2	2.02	0.42
1:O:206:ASP:HB3	2:P:241:ASN:OD1	2.20	0.42
1:O:249:GLY:HA2	1:O:259:PHE:HA	2.00	0.42
2:D:125:ASP:OD1	2:D:176:ARG:NH2	2.50	0.42
2:F:287:ASN:HB2	2:F:328:ASP:OD1	2.20	0.42
1:G:99:PRO:CB	1:G:116:VAL:HG22	2.48	0.42
2:H:343:CYS:HA	2:H:346:VAL:HG12	2.02	0.42
2:N:115:PRO:O	2:N:118:MET:HE2	2.20	0.42
1:I:172:HIS:CE1	1:I:185:LEU:HD11	2.55	0.42
2:N:81:ILE:H	2:N:81:ILE:HG12	1.60	0.42
2:N:315:VAL:HG22	2:N:342:ILE:HD13	2.02	0.42
1:B:281:LEU:O	1:B:285:VAL:HG23	2.20	0.42
2:C:62:LYS:HE3	2:C:62:LYS:HB3	1.74	0.42
1:E:88:ARG:NH2	1:E:89:LYS:HE2	2.35	0.42
1:I:120:GLU:O	1:I:226:ASN:HA	2.19	0.42
1:K:126:TYR:OH	2:L:184:LYS:HE2	2.19	0.42
2:L:19:MET:HE1	2:L:34:LYS:CE	2.50	0.42
1:M:152:ILE:HD12	1:M:152:ILE:HA	1.81	0.42
1:O:35:GLN:HE21	1:O:35:GLN:HB2	1.51	0.42
1:B:39:ARG:CD	1:B:60:SER:CB	2.98	0.42
1:B:120:GLU:OE1	1:B:122:THR:OG1	2.38	0.42
1:G:276:ASN:ND2	1:G:315:THR:OG1	2.53	0.42
2:H:293:LEU:HD23	2:H:293:LEU:HA	1.93	0.42
1:I:41:VAL:CG2	1:I:87:LEU:HD21	2.48	0.42
1:O:108:THR:C	1:O:110:TYR:N	2.78	0.42
1:O:154:GLU:O	1:O:158:GLU:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:ALA:O	1:A:26:ILE:HG13	2.20	0.41
2:D:260:GLU:HG3	2:D:262:TYR:CE2	2.54	0.41
1:E:204:TYR:CD2	2:F:277:PHE:HE2	2.38	0.41
2:H:129:ILE:O	2:H:234:VAL:HA	2.19	0.41
2:J:171:ALA:HA	2:J:176:ARG:HB2	2.02	0.41
1:K:224:MET:CE	1:K:229:GLY:N	2.83	0.41
2:L:288:PRO:HG2	2:L:339:THR:HG22	2.02	0.41
1:O:29:ALA:HB1	1:O:336:LYS:HE3	2.01	0.41
1:E:21:ALA:HA	1:E:24:MET:HE3	2.00	0.41
1:E:195:CYS:C	1:E:197:ASP:H	2.27	0.41
1:G:155:PHE:CE1	1:G:251:ILE:HG13	2.55	0.41
1:G:162:ASN:OD1	1:G:162:ASN:N	2.53	0.41
1:M:231:ILE:HD13	2:N:242:ILE:HD13	2.01	0.41
2:N:136:GLU:HG3	2:N:153:LYS:HB2	2.02	0.41
1:O:126:TYR:CD1	2:P:187:ILE:HG21	2.55	0.41
2:P:19:MET:C	2:P:20:LEU:HD12	2.44	0.41
2:C:40:ALA:HB3	2:C:42:VAL:HG23	2.02	0.41
2:D:178:LYS:HD2	2:D:212:GLU:HG3	2.03	0.41
2:F:298:MET:HE2	2:F:298:MET:HB3	1.96	0.41
1:I:235:LEU:CD1	2:J:224:VAL:HG21	2.51	0.41
2:L:279:GLN:NE2	2:L:287:ASN:H	2.04	0.41
2:P:270:GLU:CD	2:P:274:ARG:HH11	2.27	0.41
2:D:105:PHE:CE2	2:D:162:ARG:HG2	2.55	0.41
1:E:9:LEU:O	1:E:38:GLU:HA	2.20	0.41
1:E:165:ARG:HD3	1:E:165:ARG:HA	1.67	0.41
2:F:19:MET:HE3	2:F:46:PHE:CD1	2.55	0.41
2:H:37:PHE:CE2	2:H:42:VAL:HG11	2.56	0.41
2:H:134:GLU:OE1	2:H:159:LYS:NZ	2.52	0.41
2:H:288:PRO:O	2:H:292:LEU:HG	2.20	0.41
2:L:180:THR:HG23	2:L:212:GLU:HB2	2.03	0.41
1:M:39:ARG:HE	1:M:56:GLU:C	2.27	0.41
1:M:206:ASP:OD1	1:M:206:ASP:N	2.51	0.41
2:N:81:ILE:O	2:N:81:ILE:HG22	2.19	0.41
1:O:105:GLY:HA3	1:O:308:ILE:HG22	2.03	0.41
1:O:287:MET:CG	1:O:291:MET:HE2	2.50	0.41
2:P:116:GLY:HA3	2:P:320:LYS:HA	2.02	0.41
1:B:73:LYS:HE3	1:B:73:LYS:HB3	1.77	0.41
2:C:81:ILE:HD12	2:C:95:ASP:CA	2.40	0.41
2:D:182:VAL:HA	2:D:214:MET:O	2.21	0.41
1:E:80:HIS:HE1	1:E:83:MET:HE1	1.85	0.41
1:I:99:PRO:HB3	1:I:116:VAL:HG22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:129:ILE:HD13	2:P:142:HIS:CD2	2.56	0.41
1:O:149:SER:HA	1:O:152:ILE:HG22	2.02	0.41
1:O:223:VAL:C	1:O:224:MET:HG3	2.45	0.41
2:P:16:PRO:HB2	2:P:73:ASN:CG	2.45	0.41
2:P:36:VAL:HG11	2:P:292:LEU:HD22	2.02	0.41
2:P:182:VAL:HA	2:P:214:MET:O	2.20	0.41
1:B:232:LEU:HD12	1:B:232:LEU:HA	1.79	0.41
2:H:42:VAL:HG12	2:H:44:VAL:HG13	2.03	0.41
2:J:343:CYS:HA	2:J:346:VAL:HG12	2.02	0.41
2:P:168:PHE:O	2:P:172:THR:HG23	2.20	0.41
1:A:123:GLU:HB2	1:A:143:LEU:O	2.21	0.41
2:D:168:PHE:HB3	2:D:206:TYR:CD2	2.56	0.41
2:F:130:ARG:HG3	2:F:240:GLY:HA3	2.03	0.41
1:G:283:SER:O	1:G:286:MET:HB2	2.21	0.41
1:M:126:TYR:CE1	1:M:227:LEU:HD12	2.56	0.41
1:A:94:TYR:CE2	1:A:151:ARG:HG2	2.56	0.41
1:A:224:MET:HE2	1:A:224:MET:HB3	1.96	0.41
2:D:96:MET:HE2	2:D:99:ARG:HD3	2.00	0.41
2:D:165:LYS:HA	2:D:202:VAL:HG11	2.03	0.41
2:L:184:LYS:HD2	2:L:215:ILE:CG2	2.51	0.41
1:O:10:ILE:HG21	1:O:41:VAL:HG22	2.02	0.41
2:C:127:VAL:HG11	2:C:170:TYR:CE1	2.56	0.41
2:C:183:HIS:CE1	2:C:196:LEU:HD11	2.56	0.41
1:A:4:VAL:HA	1:A:33:PRO:HB2	2.03	0.41
1:A:197:ASP:HB3	1:K:162:ASN:OD1	2.21	0.41
2:D:325:ARG:HB2	2:D:331:GLY:HA3	2.03	0.41
1:E:126:TYR:N	1:E:126:TYR:CD1	2.88	0.41
1:E:222:LEU:HB3	1:E:224:MET:HE1	2.02	0.41
1:E:224:MET:HE2	1:E:224:MET:HB3	1.74	0.41
1:G:14:GLY:C	1:G:16:GLY:H	2.29	0.41
1:G:65:LYS:O	1:G:66:MET:HE2	2.20	0.41
2:H:287:ASN:ND2	2:H:329:MET:HG3	2.35	0.41
2:H:300:ARG:NH2	2:H:305:GLU:OE1	2.54	0.41
1:I:203:MET:HE1	1:I:211:ASN:HD22	1.86	0.41
2:J:165:LYS:HE2	2:J:169:ASP:OD2	2.21	0.41
2:J:197:GLN:NE2	2:J:201:GLU:OE1	2.54	0.41
2:J:200:GLU:HG3	2:J:211:PHE:HE2	1.85	0.41
2:J:343:CYS:O	2:J:346:VAL:HG12	2.20	0.41
1:K:20:SER:O	1:K:24:MET:HG3	2.21	0.41
1:K:287:MET:HE2	1:K:287:MET:HB3	1.80	0.41
2:L:235:MET:HB2	2:L:235:MET:HE3	1.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:288:PRO:CG	2:L:339:THR:HG22	2.51	0.41
2:L:311:ILE:O	2:L:315:VAL:HG23	2.21	0.41
1:M:68:LEU:HD23	1:M:68:LEU:HA	1.76	0.41
1:O:332:CYS:C	1:O:335:VAL:HG12	2.46	0.41
2:P:324:VAL:HG21	2:P:341:GLU:HG2	2.03	0.41
2:P:324:VAL:HG11	2:P:338:PHE:HA	2.03	0.41
1:A:317:ASP:OD1	1:A:317:ASP:N	2.51	0.41
2:D:107:ASN:HB2	2:D:132:GLN:CD	2.46	0.41
1:G:291:MET:HE3	1:G:293:LEU:HD12	2.04	0.41
2:J:160:SER:O	2:J:163:ILE:HG22	2.21	0.41
1:K:157:PHE:O	1:K:161:ARG:HG3	2.20	0.41
1:K:171:VAL:HB	1:K:224:MET:HG2	2.01	0.41
2:N:143:GLU:HG2	2:N:145:ALA:O	2.20	0.41
1:O:118:ILE:CD1	1:O:156:ALA:HA	2.51	0.41
2:P:25:VAL:HG21	2:P:286:ALA:CB	2.51	0.41
1:B:185:LEU:HD23	1:B:185:LEU:HA	1.77	0.40
2:F:20:LEU:HD21	2:F:66:VAL:HG21	2.03	0.40
2:F:81:ILE:HG21	2:F:95:ASP:CA	2.50	0.40
2:H:19:MET:HE2	2:H:30:MET:SD	2.61	0.40
2:H:172:THR:HG22	2:H:209:ILE:HD11	2.03	0.40
2:H:325:ARG:O	2:H:331:GLY:HA3	2.22	0.40
1:K:119:ARG:HD2	1:K:229:GLY:C	2.46	0.40
1:M:268:ASP:OD1	1:M:268:ASP:N	2.52	0.40
1:G:69:LYS:O	1:G:260:GLU:HA	2.21	0.40
1:G:179:MET:HB3	2:H:149:ILE:HG21	2.03	0.40
2:J:293:LEU:HD23	2:J:293:LEU:HA	1.75	0.40
1:O:289:ARG:HG2	1:O:289:ARG:NH1	2.36	0.40
2:P:119:THR:O	2:P:122:ASN:ND2	2.53	0.40
2:P:188:MET:O	2:P:192:ASP:HB2	2.21	0.40
1:A:118:ILE:HD12	1:A:221:VAL:HG13	2.04	0.40
1:E:109:PRO:HB3	2:F:120:ARG:HB3	2.04	0.40
1:E:150:LYS:HB2	1:E:187:LYS:HD3	2.03	0.40
2:F:336:HIS:CE1	2:J:336:HIS:CE1	3.09	0.40
1:I:185:LEU:HA	1:I:185:LEU:HD23	1.68	0.40
2:N:170:TYR:CZ	2:N:174:LYS:HG3	2.57	0.40
2:P:345:ARG:O	2:P:345:ARG:HG3	2.21	0.40
1:B:119:ARG:HD2	1:B:230:ASP:N	2.37	0.40
2:C:178:LYS:HD3	2:C:210:LYS:HD3	2.04	0.40
1:A:214:GLN:O	2:D:120:ARG:NH2	2.47	0.40
2:D:129:ILE:CD1	2:D:167:ALA:HA	2.52	0.40
1:G:184:PHE:CE1	1:G:225:PRO:HD3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:130:GLU:OE1	2:J:190:LEU:N	2.54	0.40
1:O:184:PHE:CE2	1:O:225:PRO:HD3	2.56	0.40
2:P:65:GLN:O	2:P:68:SER:OG	2.40	0.40
2:P:131:GLU:O	2:P:237:ASN:HA	2.22	0.40
2:P:220:CYS:O	2:P:224:VAL:HG23	2.21	0.40
1:A:19:ILE:O	1:A:23:VAL:HG23	2.21	0.40
1:A:276:ASN:HA	1:A:277:PRO:HD3	1.89	0.40
2:D:189:LYS:O	2:D:193:GLY:HA3	2.21	0.40
1:E:119:ARG:HD2	1:E:229:GLY:C	2.46	0.40
1:E:143:LEU:HD12	1:E:143:LEU:HA	1.87	0.40
2:H:110:HIS:HE2	2:H:260:GLU:CD	2.29	0.40
1:I:83:MET:HE3	1:I:83:MET:HB3	1.91	0.40
1:I:125:GLU:HG2	1:I:142:LYS:HG3	2.04	0.40
1:K:232:LEU:HD12	1:K:232:LEU:HA	1.75	0.40
2:P:178:LYS:O	2:P:231:ASP:HB3	2.22	0.40
2:P:217:ASP:N	2:P:217:ASP:OD1	2.55	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)
2:C:350:ASP:OD1	2:F:344:ARG:NH1[3_545]	2.03	0.17

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	333/339 (98%)	319 (96%)	14 (4%)	0	100	100
1	B	329/339 (97%)	308 (94%)	21 (6%)	0	100	100
1	E	335/339 (99%)	322 (96%)	13 (4%)	0	100	100
1	G	333/339 (98%)	310 (93%)	23 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	I	329/339 (97%)	312 (95%)	17 (5%)	0	100	100
1	K	329/339 (97%)	310 (94%)	19 (6%)	0	100	100
1	M	333/339 (98%)	306 (92%)	26 (8%)	1 (0%)	36	55
1	O	333/339 (98%)	300 (90%)	33 (10%)	0	100	100
2	C	313/352 (89%)	294 (94%)	19 (6%)	0	100	100
2	D	313/352 (89%)	295 (94%)	18 (6%)	0	100	100
2	F	309/352 (88%)	289 (94%)	20 (6%)	0	100	100
2	H	313/352 (89%)	296 (95%)	17 (5%)	0	100	100
2	J	311/352 (88%)	286 (92%)	25 (8%)	0	100	100
2	L	311/352 (88%)	296 (95%)	15 (5%)	0	100	100
2	N	310/352 (88%)	293 (94%)	17 (6%)	0	100	100
2	P	314/352 (89%)	295 (94%)	19 (6%)	0	100	100
All	All	5148/5528 (93%)	4831 (94%)	316 (6%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	109	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	266/276 (96%)	266 (100%)	0	100	100
1	B	261/276 (95%)	260 (100%)	1 (0%)	84	92
1	E	266/276 (96%)	266 (100%)	0	100	100
1	G	261/276 (95%)	260 (100%)	1 (0%)	84	92
1	I	259/276 (94%)	259 (100%)	0	100	100
1	K	260/276 (94%)	260 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	M	263/276 (95%)	263 (100%)	0	100	100
1	O	263/276 (95%)	257 (98%)	6 (2%)	44	68
2	C	252/296 (85%)	251 (100%)	1 (0%)	84	92
2	D	248/296 (84%)	248 (100%)	0	100	100
2	F	249/296 (84%)	249 (100%)	0	100	100
2	H	249/296 (84%)	249 (100%)	0	100	100
2	J	253/296 (86%)	252 (100%)	1 (0%)	84	92
2	L	248/296 (84%)	248 (100%)	0	100	100
2	N	252/296 (85%)	252 (100%)	0	100	100
2	P	248/296 (84%)	248 (100%)	0	100	100
All	All	4098/4576 (90%)	4088 (100%)	10 (0%)	87	95

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	336	LYS
2	C	323	LYS
1	G	104	GLU
2	J	61	GLU
1	O	35	GLN
1	O	269	ILE
1	O	276	ASN
1	O	280	LEU
1	O	296	HIS
1	O	314	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	254	ASN
2	C	229	GLN
2	C	301	HIS
2	C	336	HIS
1	A	5	GLN
1	A	290	HIS
2	D	47	GLN
2	D	107	ASN

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Mol	Chain	Res	Type
2	D	307	HIS
1	E	163	ASN
1	E	186	GLN
1	E	214	GLN
2	F	49	HIS
2	F	121	HIS
2	F	161	GLN
1	G	163	ASN
1	G	218	GLN
1	G	250	ASN
1	G	296	HIS
2	H	107	ASN
2	H	226	ASN
2	H	245	ASN
1	I	96	ASN
1	I	163	ASN
1	I	218	GLN
1	I	250	ASN
2	J	50	HIS
2	J	218	ASN
2	J	222	GLN
2	J	226	ASN
2	J	245	ASN
2	J	336	HIS
1	K	40	ASN
1	K	163	ASN
1	K	167	ASN
1	K	218	GLN
1	K	276	ASN
1	K	290	HIS
2	L	47	GLN
2	L	107	ASN
2	L	218	ASN
2	L	222	GLN
2	L	229	GLN
2	L	279	GLN
1	M	35	GLN
1	M	296	HIS
2	N	279	GLN
1	O	35	GLN
1	O	218	GLN
1	O	290	HIS

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Mol	Chain	Res	Type
2	P	47	GLN
2	P	50	HIS
2	P	107	ASN
2	P	132	GLN
2	P	225	GLN
2	P	307	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 [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	335/339 (98%)	-0.27	1 (0%) 90 89	31, 46, 77, 91	0
1	B	333/339 (98%)	-0.14	4 (1%) 76 71	34, 50, 73, 112	0
1	E	337/339 (99%)	-0.29	3 (0%) 81 76	34, 45, 67, 95	0
1	G	335/339 (98%)	0.05	3 (0%) 81 76	39, 55, 80, 94	0
1	I	333/339 (98%)	-0.25	4 (1%) 76 71	33, 46, 71, 96	0
1	K	333/339 (98%)	-0.24	2 (0%) 85 83	28, 45, 71, 87	0
1	M	335/339 (98%)	0.08	6 (1%) 67 60	32, 51, 83, 97	0
1	O	335/339 (98%)	0.29	10 (2%) 52 43	32, 59, 89, 101	0
2	C	319/352 (90%)	-0.19	8 (2%) 58 49	32, 47, 82, 98	0
2	D	319/352 (90%)	-0.11	7 (2%) 62 53	34, 49, 78, 96	0
2	F	315/352 (89%)	-0.06	12 (3%) 44 35	38, 51, 86, 102	0
2	H	319/352 (90%)	-0.03	8 (2%) 58 49	40, 52, 83, 99	0
2	J	317/352 (90%)	-0.19	7 (2%) 62 53	34, 47, 79, 98	0
2	L	317/352 (90%)	-0.12	8 (2%) 58 49	32, 48, 84, 121	0
2	N	316/352 (89%)	-0.20	4 (1%) 75 69	34, 48, 76, 104	0
2	P	320/352 (90%)	-0.06	13 (4%) 41 33	36, 50, 80, 104	0
All	All	5218/5528 (94%)	-0.11	100 (1%) 66 59	28, 50, 80, 121	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	P	92	ALA	5.3
1	B	47	PRO	5.3
2	N	63	LEU	4.7
2	F	51	LEU	4.5
1	I	46	GLY	4.2

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Mol	Chain	Res	Type	RSRZ
2	H	92	ALA	3.9
1	O	4	VAL	3.7
1	B	2	GLY	3.5
1	O	292	GLY	3.5
2	F	14	SER	3.5
1	I	50	LYS	3.4
2	D	62	LYS	3.2
1	K	47	PRO	3.2
2	F	63	LEU	3.2
2	J	277	PHE	3.2
2	D	92	ALA	3.1
2	H	62	LYS	3.1
2	J	281	VAL	3.1
1	E	2	GLY	3.1
2	P	116	GLY	3.1
2	C	52	SER	3.0
2	L	81	ILE	2.9
2	F	281	VAL	2.9
2	C	94	TYR	2.8
2	P	61	GLU	2.8
2	L	15	PHE	2.8
2	D	93	SER	2.8
2	F	93	SER	2.8
2	J	93	SER	2.8
2	H	50	HIS	2.7
2	P	50	HIS	2.7
1	E	1	THR	2.7
1	I	76	ILE	2.7
1	G	273	ASP	2.7
1	M	295	ASP	2.7
2	C	63	LEU	2.7
2	J	63	LEU	2.7
1	K	50	LYS	2.7
1	O	156	ALA	2.6
2	C	277	PHE	2.6
1	M	29	ALA	2.6
1	M	3	GLY	2.6
2	F	80	LYS	2.6
1	O	337	ASP	2.6
2	F	280	ALA	2.6
1	O	107	LYS	2.5
2	C	281	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
2	F	23	ASP	2.5
1	I	2	GLY	2.5
2	L	96	MET	2.5
2	L	349	LEU	2.5
1	G	321	ASN	2.5
2	L	52	SER	2.4
2	L	50	HIS	2.4
2	C	280	ALA	2.4
1	O	6	THR	2.4
2	P	52	SER	2.4
1	M	4	VAL	2.3
2	P	265	GLU	2.3
1	O	293	LEU	2.3
2	P	97	ARG	2.3
1	G	313	SER	2.3
2	C	81	ILE	2.3
2	H	332	TYR	2.3
2	J	115	PRO	2.3
2	P	24	GLY	2.3
2	L	92	ALA	2.3
2	H	51	LEU	2.3
2	C	93	SER	2.3
2	H	93	SER	2.3
1	A	319	GLY	2.3
2	F	92	ALA	2.3
2	D	52	SER	2.3
2	D	63	LEU	2.2
2	P	51	LEU	2.2
2	N	62	LYS	2.2
2	P	62	LYS	2.2
2	P	93	SER	2.2
2	D	95	ASP	2.2
2	P	95	ASP	2.2
2	F	96	MET	2.2
1	B	46	GLY	2.2
2	H	94	TYR	2.2
1	B	39	ARG	2.1
2	N	23	ASP	2.1
1	M	105	GLY	2.1
2	L	330	GLY	2.1
1	O	298	ALA	2.1
2	F	95	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
2	J	52	SER	2.1
2	P	81	ILE	2.1
1	O	321	ASN	2.1
2	H	95	ASP	2.1
2	J	282	GLY	2.1
1	M	57	ALA	2.0
1	E	313	SER	2.0
1	O	296	HIS	2.0
2	F	81	ILE	2.0
2	N	96	MET	2.0
2	D	94	TYR	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.