



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 11:53 AM EDT

PDB ID : 6BPE / pdb_00006bpe
Title : Plasmodium vivax reticulocyte binding protein 2b (PvRBP2b) bound to monoclonal antibody 6H1
Authors : Gruszczyk, J.; Chan, L.J.; Tham, W.H.
Deposited on : 2017-11-22
Resolution : 3.34 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

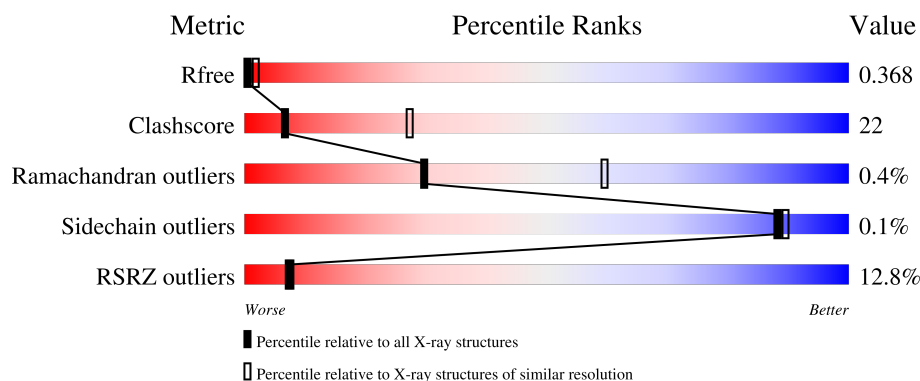
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.34 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	1434 (3.38-3.30)
Clashscore	190562	1479 (3.38-3.30)
Ramachandran outliers	187476	1456 (3.38-3.30)
Sidechain outliers	187428	1455 (3.38-3.30)
RSRZ outliers	180081	1434 (3.38-3.30)

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	307	10% 65% 32% .
1	D	307	10% 65% 33% .
1	G	307	13% 61% 38% .
1	J	307	16% 52% 44% .
2	B	254	9% 44% 38% 18%

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Mol	Chain	Length	Quality of chain
2	E	254	
2	H	254	
3	C	238	
3	F	238	
3	I	238	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17777 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 Reticulocyte binding protein 2, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2518	1612	430	467	9			
1	D	302	Total	C	N	O	S	0	0	0
			2544	1627	436	472	9			
1	G	302	Total	C	N	O	S	0	0	0
			2543	1627	436	471	9			
1	J	297	Total	C	N	O	S	0	0	0
			2500	1602	428	461	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	GLY	-	expression tag	UNP A5K736
A	165	ALA	-	expression tag	UNP A5K736
A	166	MET	-	expression tag	UNP A5K736
A	167	GLY	-	expression tag	UNP A5K736
A	168	SER	-	expression tag	UNP A5K736
D	164	GLY	-	expression tag	UNP A5K736
D	165	ALA	-	expression tag	UNP A5K736
D	166	MET	-	expression tag	UNP A5K736
D	167	GLY	-	expression tag	UNP A5K736
D	168	SER	-	expression tag	UNP A5K736
G	164	GLY	-	expression tag	UNP A5K736
G	165	ALA	-	expression tag	UNP A5K736
G	166	MET	-	expression tag	UNP A5K736
G	167	GLY	-	expression tag	UNP A5K736
G	168	SER	-	expression tag	UNP A5K736
J	164	GLY	-	expression tag	UNP A5K736
J	165	ALA	-	expression tag	UNP A5K736
J	166	MET	-	expression tag	UNP A5K736
J	167	GLY	-	expression tag	UNP A5K736
J	168	SER	-	expression tag	UNP A5K736

- Molecule 2 is a protein called Monoclonal antibody 6H1 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	209	Total	C	N	O	S	0	0	0
			1594	1013	252	321	8			
2	E	208	Total	C	N	O	S	0	0	0
			1590	1011	254	318	7			
2	H	113	Total	C	N	O	S	0	0	0
			885	560	140	180	5			

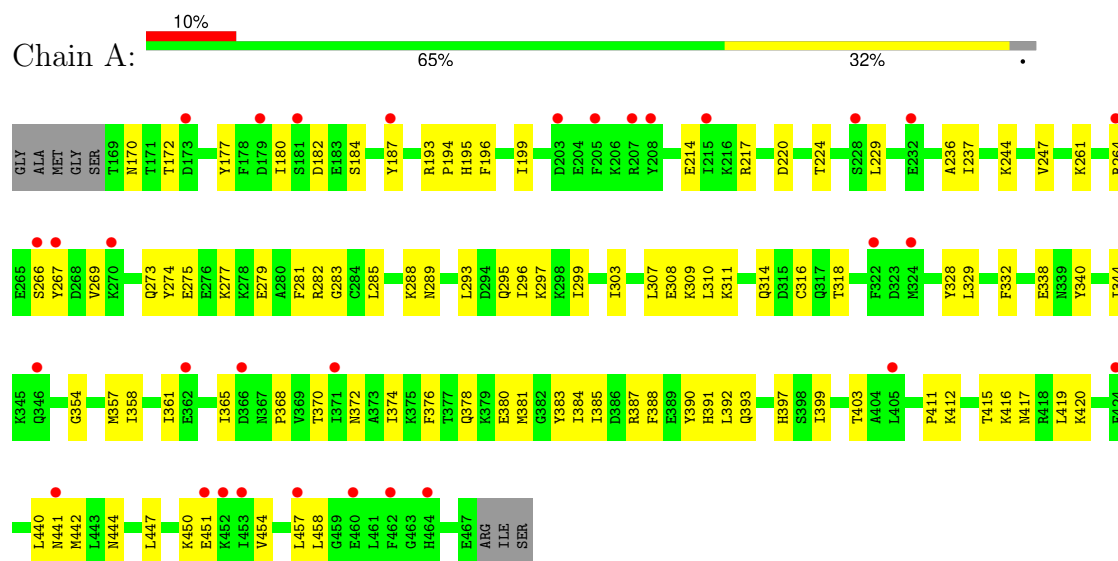
- Molecule 3 is a protein called Monoclonal antibody 6H1 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	199	Total	C	N	O	S	0	0	0
			1539	960	258	314	7			
3	F	198	Total	C	N	O	S	0	0	0
			1535	962	255	311	7			
3	I	67	Total	C	N	O	S	0	0	0
			529	334	88	105	2			

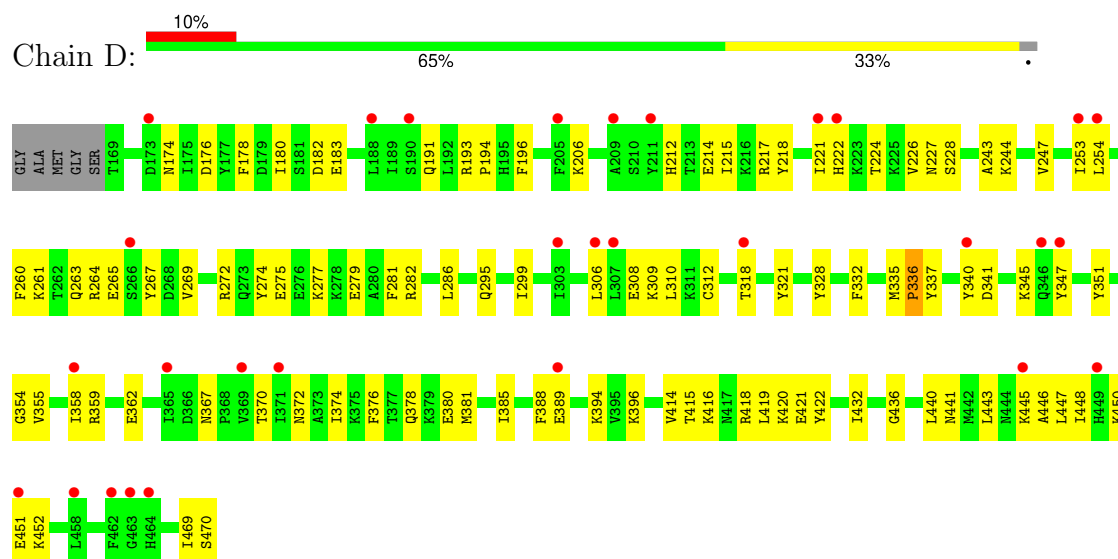
3 Residue-property plots

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.

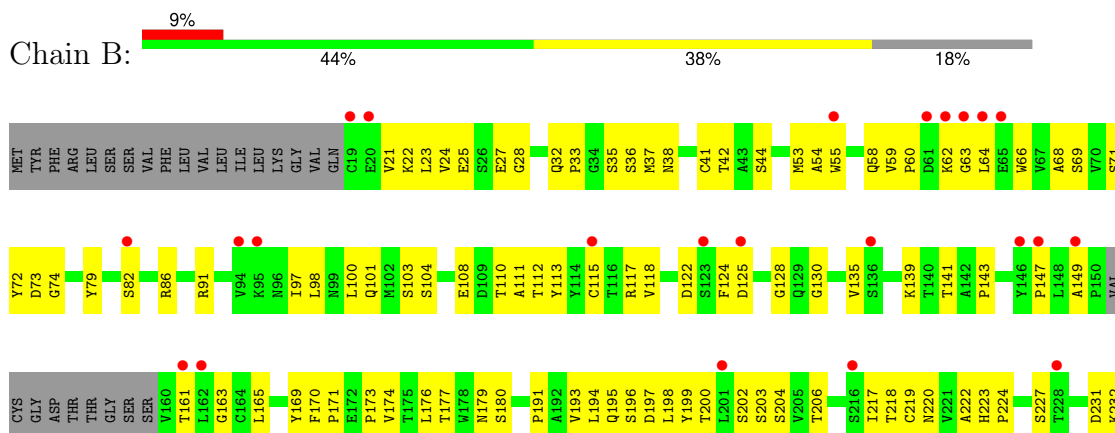
- Molecule 1: Reticulocyte binding protein 2, putative

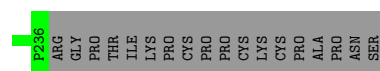


- Molecule 1: Reticulocyte binding protein 2, putative

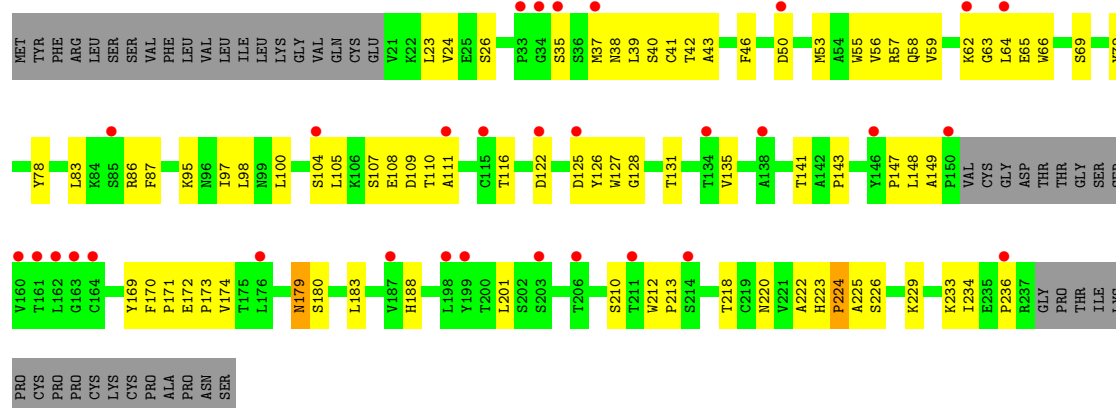


- Molecule 1: Reticulocyte binding protein 2, putative

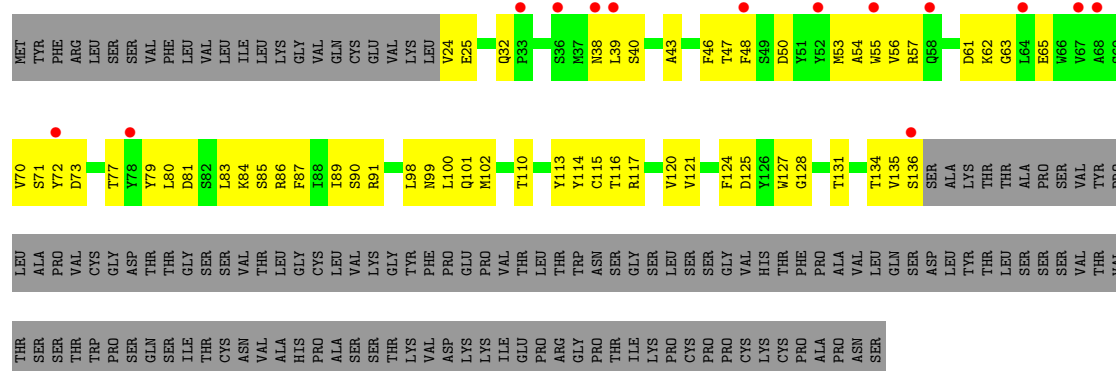




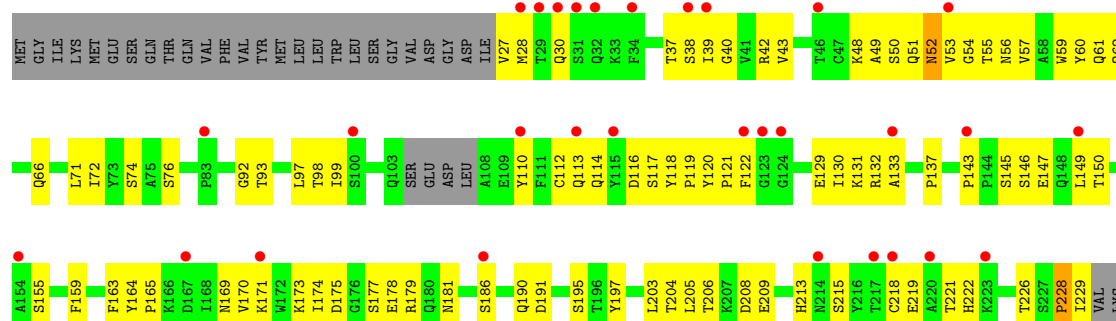
• Molecule 2: Monoclonal antibody 6H1 Fab heavy chain



• Molecule 2: Monoclonal antibody 6H1 Fab heavy chain



• Molecule 3: Monoclonal antibody 6H1 Fab light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.60Å 123.48Å 143.28Å 90.00° 105.25° 90.00°	Depositor
Resolution (Å)	73.75 – 3.34 73.75 – 3.34	Depositor EDS
% Data completeness (in resolution range)	99.2 (73.75-3.34) 99.4 (73.75-3.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 3.33Å)	Xtrriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.318 , 0.367 0.318 , 0.368	Depositor DCC
R_{free} test set	1068 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	75.2	Xtrriage
Anisotropy	0.681	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 97.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	17777	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.93 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2039e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.11	0/2567	0.27	0/3444
1	D	0.12	0/2593	0.30	0/3477
1	G	0.16	0/2592	0.36	0/3477
1	J	0.15	0/2549	0.39	0/3420
2	B	0.17	0/1632	0.46	0/2231
2	E	0.17	0/1628	0.44	0/2225
2	H	0.17	0/904	0.42	0/1229
3	C	0.15	0/1574	0.42	0/2136
3	F	0.18	0/1570	0.42	0/2128
3	I	0.29	0/541	0.78	2/732 (0.3%)
All	All	0.16	0/18150	0.40	2/24499 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	112	CYS	N-CA-C	8.66	123.18	108.02
3	I	59	TRP	N-CA-C	-5.22	99.90	108.41

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	2518	0	2523	75	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2544	0	2552	71	0
1	G	2543	0	2552	106	0
1	J	2500	0	2511	108	0
2	B	1594	0	1555	87	0
2	E	1590	0	1557	78	0
2	H	885	0	843	61	0
3	C	1539	0	1465	75	0
3	F	1535	0	1466	69	0
3	I	529	0	491	69	0
All	All	17777	0	17515	760	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 22.

All (760) 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)
3:I:57:VAL:CG2	3:I:59:TRP:CH2	1.76	1.62
3:I:57:VAL:HG21	3:I:59:TRP:CH2	1.37	1.49
3:I:57:VAL:HG23	3:I:59:TRP:CZ3	1.48	1.44
3:I:53:VAL:HG21	3:I:116:ASP:CG	1.48	1.37
2:E:56:VAL:HG23	2:E:65:GLU:O	1.24	1.30
3:I:53:VAL:HG21	3:I:116:ASP:OD1	1.33	1.29
3:I:59:TRP:CD1	3:I:112:CYS:HB3	1.67	1.28
3:I:57:VAL:HG21	3:I:59:TRP:CZ2	1.67	1.27
3:C:219:GLU:OE2	3:C:228:PRO:HB2	1.25	1.26
3:C:219:GLU:OE2	3:C:228:PRO:CB	1.83	1.25
2:E:56:VAL:CG2	2:E:65:GLU:O	1.93	1.15
3:I:57:VAL:HG23	3:I:59:TRP:CH2	1.56	1.12
3:I:53:VAL:CG2	3:I:116:ASP:CG	2.33	1.01
3:I:52:ASN:ND2	3:I:54:GLY:O	1.92	1.01
3:I:59:TRP:CD1	3:I:112:CYS:CB	2.44	1.00
3:F:113:GLN:HG3	3:F:122:PHE:HD2	1.33	0.94
3:C:28:MET:HB2	3:C:50:SER:H	1.32	0.92
2:B:62:LYS:HB3	2:B:63:GLY:HA2	1.53	0.91
2:H:114:TYR:HH	3:I:67:SER:N	1.67	0.91
1:G:311:LYS:CB	3:I:53:VAL:O	2.20	0.90
3:I:58:ALA:H	3:I:115:TYR:HE2	1.16	0.90
2:E:64:LEU:O	3:F:122:PHE:CE1	2.09	0.90
2:E:58:GLN:HE21	2:E:62:LYS:HG3	1.37	0.89
3:C:53:VAL:HG11	3:C:114:GLN:HG3	1.56	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:179:ASN:HD21	2:E:183:LEU:HD13	1.39	0.88
3:I:59:TRP:HD1	3:I:112:CYS:HB3	1.32	0.87
3:I:57:VAL:CG2	3:I:59:TRP:HH2	1.77	0.87
3:I:59:TRP:HZ3	3:I:75:ALA:HB2	1.39	0.86
3:I:57:VAL:CG2	3:I:59:TRP:CZ2	2.42	0.84
1:G:311:LYS:HB3	3:I:53:VAL:C	2.01	0.84
2:B:170:PHE:HB2	2:B:198:LEU:HD21	1.58	0.83
3:C:219:GLU:CD	3:C:228:PRO:CB	2.51	0.83
2:B:177:THR:HB	2:B:220:ASN:HB2	1.60	0.83
1:D:418:ARG:NH2	1:D:421:GLU:OE1	2.11	0.82
1:J:318:THR:H	1:J:412:LYS:HG3	1.45	0.81
1:J:205:PHE:HA	1:J:208:TYR:HD2	1.46	0.81
1:G:311:LYS:HB3	3:I:53:VAL:O	1.79	0.80
2:B:179:ASN:HA	2:B:217:ILE:HG21	1.64	0.80
3:C:137:PRO:HD3	3:C:222:HIS:HD2	1.47	0.79
3:I:55:THR:HG21	3:I:91:SER:HA	1.65	0.79
2:E:110:THR:HG23	2:E:135:VAL:H	1.47	0.79
3:F:137:PRO:HB3	3:F:163:PHE:HB3	1.64	0.78
2:E:86:ARG:NH2	2:E:109:ASP:OD2	2.17	0.77
2:H:84:LYS:NZ	2:H:85:SER:H	1.82	0.77
1:J:290:ARG:HD2	1:J:293:LEU:HD21	1.66	0.77
1:J:421:GLU:O	1:J:425:ASN:N	2.18	0.77
1:D:247:VAL:HG22	1:D:277:LYS:HB3	1.66	0.76
3:C:219:GLU:CD	3:C:228:PRO:HB2	2.10	0.76
3:I:59:TRP:CZ2	3:I:95:PHE:HB3	2.20	0.76
2:E:64:LEU:C	3:F:122:PHE:CZ	2.37	0.75
3:C:206:THR:OG1	3:C:208:ASP:OD1	2.04	0.75
1:J:214:GLU:HG3	1:J:311:LYS:H	1.50	0.75
1:G:224:THR:HG21	2:H:73:ASP:OD2	1.87	0.75
3:F:172:TRP:HE1	3:F:201:SER:HB3	1.50	0.74
3:C:205:LEU:HD11	3:C:209:GLU:HB2	1.66	0.74
1:D:367:ASN:OD1	1:D:370:THR:OG1	2.05	0.74
2:E:58:GLN:NE2	2:E:59:VAL:O	2.20	0.74
2:E:180:SER:H	2:E:220:ASN:HD21	1.35	0.74
2:B:27:GLU:OE2	2:B:27:GLU:N	2.21	0.74
3:F:209:GLU:HA	3:F:212:ARG:HD3	1.70	0.73
3:C:129:GLU:OE2	3:C:197:TYR:OH	2.07	0.73
2:H:55:TRP:NE1	2:H:115:CYS:SG	2.62	0.73
1:A:217:ARG:NH2	2:B:72:TYR:OH	2.20	0.73
2:E:147:PRO:O	3:F:145:SER:OG	2.06	0.72
2:B:117:ARG:NE	2:B:125:ASP:OD2	2.20	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:110:THR:HG22	2:B:135:VAL:H	1.54	0.72
3:C:38:SER:HB3	3:C:131:LYS:HB2	1.72	0.72
3:I:43:VAL:HG23	3:I:98:THR:O	1.91	0.71
1:J:217:ARG:HH22	1:J:309:LYS:HD3	1.53	0.71
3:I:53:VAL:CG2	3:I:116:ASP:OD1	2.27	0.71
1:A:172:THR:HG21	1:A:450:LYS:HE2	1.73	0.70
3:F:174:ILE:O	3:F:174:ILE:HD12	1.91	0.70
1:G:311:LYS:HB2	3:I:53:VAL:O	1.91	0.70
3:C:219:GLU:OE2	3:C:228:PRO:CG	2.38	0.70
2:E:43:ALA:HB1	2:E:46:PHE:HE1	1.57	0.70
2:B:82:SER:O	2:B:86:ARG:NH2	2.24	0.70
1:D:178:PHE:O	1:D:191:GLN:NE2	2.21	0.70
3:C:219:GLU:CD	3:C:228:PRO:HB3	2.17	0.69
3:F:78:ARG:NH2	3:F:84:ASP:OD1	2.26	0.69
3:I:58:ALA:HB3	3:I:115:TYR:CZ	2.28	0.69
2:B:25:GLU:HG2	2:B:55:TRP:HH2	1.55	0.69
2:B:220:ASN:HA	2:B:231:ASP:OD1	1.93	0.69
1:D:347:TYR:OH	1:D:447:LEU:O	2.10	0.68
1:G:311:LYS:HD3	3:I:53:VAL:O	1.93	0.68
3:C:28:MET:HE2	3:C:53:VAL:HG13	1.76	0.68
2:H:53:MET:HB3	2:H:98:LEU:HD22	1.76	0.68
2:B:191:PRO:HG2	3:C:186:SER:HB2	1.75	0.68
2:E:39:LEU:HD22	2:E:131:THR:HG21	1.76	0.68
1:G:267:TYR:HB3	1:G:372:ASN:HD22	1.59	0.68
1:J:344:ILE:HD11	1:J:385:ILE:HG23	1.75	0.68
1:J:454:VAL:O	1:J:458:LEU:N	2.23	0.68
1:J:205:PHE:HA	1:J:208:TYR:CD2	2.29	0.67
1:A:247:VAL:HG22	1:A:277:LYS:HB3	1.76	0.67
2:H:80:LEU:HB3	2:H:83:LEU:HB2	1.76	0.67
3:C:116:ASP:OD1	3:C:117:SER:OG	2.13	0.67
1:D:274:TYR:OH	1:D:380:GLU:OE2	2.13	0.67
3:C:59:TRP:HB2	3:C:72:ILE:HB	1.76	0.67
2:E:56:VAL:HG22	2:E:65:GLU:O	1.95	0.66
1:A:357:MET:HG3	1:J:346:GLN:HG2	1.75	0.66
3:C:28:MET:SD	3:C:51:GLN:N	2.66	0.66
1:G:215:ILE:HG23	1:G:306:LEU:HD22	1.77	0.66
1:G:267:TYR:HB3	1:G:372:ASN:ND2	2.10	0.66
3:I:58:ALA:HB3	3:I:115:TYR:OH	1.95	0.66
3:I:60:TYR:OH	3:I:113:GLN:OE1	2.13	0.66
1:J:217:ARG:NE	1:J:306:LEU:HG	2.10	0.66
2:B:115:CYS:O	2:B:128:GLY:N	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:179:ASN:HB3	2:E:218:THR:H	1.60	0.66
3:F:174:ILE:HG22	3:F:216:TYR:HD1	1.61	0.66
1:G:419:LEU:HD12	1:G:423:TYR:CE2	2.30	0.66
2:E:83:LEU:O	2:E:83:LEU:HD12	1.95	0.65
1:G:416:LYS:HA	1:G:419:LEU:HD21	1.77	0.65
2:B:108:GLU:N	2:B:108:GLU:OE1	2.29	0.65
2:E:171:PRO:O	2:E:223:HIS:NE2	2.30	0.65
1:G:358:ILE:HD11	1:G:371:ILE:HG23	1.79	0.65
1:G:254:LEU:HD12	1:G:270:LYS:HD3	1.79	0.65
2:H:84:LYS:HZ2	2:H:85:SER:H	1.43	0.65
2:B:62:LYS:CB	2:B:63:GLY:HA2	2.23	0.65
3:I:57:VAL:CG2	3:I:59:TRP:CZ3	2.31	0.65
1:D:174:ASN:ND2	1:D:176:ASP:OD1	2.27	0.64
1:J:302:GLU:O	1:J:306:LEU:HD13	1.97	0.64
1:A:370:THR:HG22	1:A:457:LEU:HD21	1.80	0.64
3:C:56:ASN:ND2	3:C:116:ASP:OD2	2.31	0.64
2:E:218:THR:HG22	2:E:233:LYS:HA	1.78	0.64
3:C:60:TYR:OH	3:C:113:GLN:OE1	2.13	0.64
3:C:190:GLN:NE2	3:C:191:ASP:O	2.30	0.64
2:E:58:GLN:O	2:E:111:ALA:HB1	1.98	0.64
2:H:80:LEU:HD23	2:H:83:LEU:HD23	1.79	0.64
3:C:62:GLN:NE2	3:C:66:GLN:O	2.31	0.63
2:B:25:GLU:HG2	2:B:55:TRP:CH2	2.32	0.63
3:F:147:GLU:N	3:F:147:GLU:OE1	2.31	0.63
2:B:41:CYS:HB3	2:B:98:LEU:HB3	1.79	0.63
2:E:147:PRO:HB3	2:E:234:ILE:HG12	1.81	0.63
3:F:172:TRP:CD1	3:F:183:VAL:HG21	2.34	0.63
1:A:309:LYS:NZ	2:B:122:ASP:OD1	2.31	0.63
2:B:58:GLN:HB3	2:B:112:THR:HB	1.79	0.63
3:F:206:THR:O	3:F:210:TYR:N	2.25	0.63
1:A:229:LEU:HD21	1:A:296:ILE:HG23	1.80	0.63
2:B:194:LEU:HA	2:B:199:TYR:HA	1.79	0.63
3:I:59:TRP:NE1	3:I:112:CYS:CB	2.62	0.63
3:C:30:GLN:HG2	3:C:49:ALA:HA	1.81	0.62
2:E:116:THR:HG22	2:E:127:TRP:HA	1.80	0.62
2:H:72:TYR:HA	2:H:91:ARG:NH1	2.14	0.62
2:E:38:ASN:HD22	2:E:39:LEU:H	1.47	0.62
2:E:55:TRP:HE1	2:E:98:LEU:HD22	1.64	0.62
3:C:114:GLN:HG2	3:C:116:ASP:H	1.64	0.62
2:E:26:SER:OG	2:E:40:SER:N	2.27	0.62
1:G:231:ASN:OD1	1:G:232:GLU:N	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:193:ARG:HG3	1:D:194:PRO:HA	1.81	0.62
2:E:86:ARG:HH21	2:E:105:LEU:HA	1.65	0.62
3:F:149:LEU:O	3:F:207:LYS:HD2	1.99	0.62
3:F:169:ASN:HB3	3:F:221:THR:HB	1.80	0.62
1:J:232:GLU:HA	1:J:235:ARG:HD3	1.81	0.62
1:D:269:VAL:HG23	1:D:272:ARG:HH21	1.65	0.62
2:E:141:THR:O	2:E:226:SER:OG	2.18	0.62
1:G:295:GLN:O	1:G:299:ILE:HD12	1.98	0.62
3:C:113:GLN:HB2	3:C:122:PHE:HD1	1.65	0.62
1:G:309:LYS:HG3	2:H:121:VAL:HG12	1.81	0.61
1:A:440:LEU:O	1:A:444:ASN:ND2	2.32	0.61
1:J:421:GLU:HA	1:J:424:PHE:HB3	1.82	0.61
1:A:318:THR:OG1	1:A:411:PRO:O	2.14	0.61
3:I:59:TRP:HZ2	3:I:95:PHE:HB3	1.66	0.61
3:I:89:SER:O	3:I:96:THR:OG1	2.19	0.61
2:E:41:CYS:HB3	2:E:98:LEU:HB3	1.83	0.61
3:F:72:ILE:HD13	3:F:78:ARG:HB3	1.83	0.61
3:F:113:GLN:HG3	3:F:122:PHE:CD2	2.25	0.61
1:G:292:ASN:O	1:G:296:ILE:N	2.32	0.61
2:H:25:GLU:HG3	2:H:115:CYS:HB2	1.82	0.61
1:J:375:LYS:HA	1:J:378:GLN:HG3	1.83	0.61
1:J:214:GLU:HG3	1:J:310:LEU:HD12	1.81	0.60
1:J:382:GLY:HA2	1:J:385:ILE:HD12	1.83	0.60
3:C:137:PRO:HB3	3:C:163:PHE:HB3	1.83	0.60
3:I:59:TRP:NE1	3:I:112:CYS:HB3	2.16	0.60
2:B:35:SER:HB2	2:B:104:SER:H	1.66	0.60
2:E:95:LYS:HB3	2:E:97:ILE:HG12	1.84	0.60
1:G:386:ASP:HA	1:G:389:GLU:HB2	1.84	0.60
2:H:89:ILE:HA	2:H:100:LEU:HA	1.83	0.60
1:J:450:LYS:HA	1:J:453:ILE:HD12	1.84	0.60
2:H:73:ASP:N	2:H:73:ASP:OD1	2.34	0.60
2:E:105:LEU:HD12	2:E:105:LEU:H	1.65	0.60
1:G:204:GLU:OE1	1:G:328:TYR:OH	2.13	0.60
2:H:39:LEU:HD13	2:H:131:THR:HG21	1.84	0.60
1:A:417:ASN:HA	1:A:420:LYS:HD2	1.82	0.60
2:E:223:HIS:HD2	2:E:224:PRO:HD2	1.65	0.60
2:H:53:MET:HB3	2:H:98:LEU:CD2	2.32	0.60
3:F:185:ASN:HB3	3:F:201:SER:HA	1.83	0.59
3:F:55:THR:HG22	3:F:55:THR:O	2.02	0.59
1:G:220:ASP:O	1:G:224:THR:OG1	2.19	0.59
3:C:27:VAL:HG21	3:C:119:PRO:HD2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:38:ASN:ND2	2:E:100:LEU:O	2.35	0.59
1:D:337:TYR:CE2	1:D:396:LYS:HB2	2.37	0.59
2:H:32:GLN:NE2	2:H:136:SER:O	2.36	0.59
2:B:54:ALA:HB2	2:B:69:SER:HB3	1.84	0.59
1:J:374:ILE:HA	1:J:457:LEU:HD13	1.85	0.59
1:A:193:ARG:HG3	1:A:194:PRO:HA	1.83	0.59
1:G:425:ASN:OD1	1:G:426:ILE:N	2.36	0.59
3:C:213:HIS:ND1	3:C:215:SER:O	2.36	0.58
2:E:38:ASN:HD22	2:E:39:LEU:N	2.00	0.58
1:G:348:LYS:O	1:G:352:LEU:HG	2.03	0.58
1:J:200:TYR:OH	1:J:438:ASP:OD2	2.20	0.58
2:B:24:VAL:HG12	2:B:42:THR:HB	1.84	0.58
2:E:24:VAL:HG22	2:E:42:THR:HB	1.85	0.58
1:J:293:LEU:HA	1:J:296:ILE:HD12	1.85	0.58
1:J:365:ILE:HD12	1:J:366:ASP:H	1.69	0.58
2:B:28:GLY:N	2:B:37:MET:HE3	2.19	0.58
2:E:37:MET:SD	2:E:38:ASN:N	2.77	0.58
2:H:62:LYS:N	2:H:63:GLY:HA2	2.19	0.58
2:B:71:SER:O	2:B:91:ARG:NH1	2.36	0.57
3:C:164:TYR:CD1	3:C:165:PRO:HA	2.39	0.57
3:F:173:LYS:HB3	3:F:178:GLU:HB3	1.86	0.57
1:J:169:THR:HG21	1:J:452:LYS:HB3	1.85	0.57
1:J:254:LEU:HD11	1:J:270:LYS:HB3	1.86	0.57
2:E:56:VAL:HG23	2:E:65:GLU:C	2.18	0.57
1:G:409:VAL:HG13	1:G:411:PRO:HD3	1.86	0.57
1:G:311:LYS:CD	3:I:53:VAL:O	2.52	0.57
1:A:374:ILE:HD12	1:A:454:VAL:HG23	1.86	0.57
3:I:59:TRP:CZ3	3:I:75:ALA:HB2	2.29	0.57
1:A:261:LYS:HZ2	1:A:264:ARG:HD2	1.69	0.57
2:E:65:GLU:N	3:F:122:PHE:CZ	2.72	0.57
2:H:117:ARG:O	2:H:125:ASP:N	2.34	0.57
1:J:415:THR:O	1:J:419:LEU:N	2.38	0.57
1:D:310:LEU:O	1:D:416:LYS:NZ	2.34	0.57
1:A:267:TYR:CZ	1:A:372:ASN:HB3	2.39	0.57
1:D:469:ILE:HD12	1:D:470:SER:N	2.19	0.57
2:E:83:LEU:HD22	2:E:87:PHE:CD1	2.40	0.57
3:F:155:SER:HA	3:F:204:THR:HA	1.87	0.57
3:F:174:ILE:HG22	3:F:216:TYR:CD1	2.39	0.57
2:B:35:SER:HB2	2:B:104:SER:N	2.19	0.57
3:F:158:CYS:HB2	3:F:172:TRP:CZ2	2.40	0.56
2:H:48:PHE:HB3	2:H:91:ARG:HH22	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:196:PHE:HB2	1:J:199:ILE:HG12	1.86	0.56
2:H:110:THR:OG1	2:H:134:THR:HA	2.04	0.56
1:J:247:VAL:O	1:J:251:ILE:HG13	2.05	0.56
2:E:43:ALA:HB1	2:E:46:PHE:CE1	2.41	0.56
2:H:87:PHE:HB3	2:H:102:MET:HA	1.86	0.56
3:I:86:PHE:CZ	3:I:98:THR:O	2.59	0.56
1:A:180:ILE:HD12	1:A:282:ARG:HB3	1.88	0.56
3:C:132:ARG:HG3	3:C:133:ALA:H	1.70	0.56
3:I:113:GLN:HE21	3:I:120:TYR:HB3	1.70	0.56
1:D:351:TYR:HD2	1:D:381:MET:HE3	1.70	0.56
1:D:355:VAL:HG13	1:D:359:ARG:HH22	1.71	0.56
3:F:116:ASP:OD1	3:F:117:SER:OG	2.23	0.56
3:I:61:GLN:HG2	3:I:109:GLU:O	2.04	0.56
3:F:215:SER:HA	3:F:232:SER:HA	1.87	0.56
1:D:370:THR:O	1:D:374:ILE:HG13	2.05	0.56
1:D:469:ILE:HD12	1:D:470:SER:H	1.70	0.56
3:F:173:LYS:HA	3:F:178:GLU:HB3	1.87	0.56
2:H:57:ARG:NH1	2:H:65:GLU:OE1	2.39	0.56
3:I:52:ASN:ND2	3:I:54:GLY:C	2.62	0.56
2:E:58:GLN:NE2	2:E:62:LYS:HG3	2.14	0.56
1:G:309:LYS:HD2	2:H:121:VAL:HA	1.86	0.56
1:J:201:PHE:CE2	1:J:431:SER:HB2	2.41	0.56
3:C:59:TRP:CD2	3:C:97:LEU:HD12	2.42	0.55
1:J:225:LYS:O	1:J:229:LEU:HD23	2.07	0.55
1:D:341:ASP:O	1:D:345:LYS:HG2	2.06	0.55
1:G:370:THR:HG21	1:G:461:LEU:HB2	1.88	0.55
2:H:24:VAL:N	2:H:25:GLU:OE1	2.39	0.55
1:A:220:ASP:O	1:A:224:THR:OG1	2.23	0.55
1:D:308:GLU:OE1	1:D:420:LYS:NZ	2.39	0.55
3:F:113:GLN:HA	3:F:122:PHE:HB3	1.88	0.55
3:F:208:ASP:OD1	3:F:208:ASP:N	2.40	0.55
2:B:147:PRO:O	3:C:145:SER:OG	2.20	0.55
2:H:110:THR:HB	2:H:135:VAL:HG22	1.88	0.55
1:A:237:ILE:HD12	1:A:285:LEU:HD23	1.88	0.55
3:C:169:ASN:HB3	3:C:221:THR:HB	1.89	0.55
1:D:440:LEU:HA	1:D:443:LEU:HD12	1.88	0.55
1:A:332:PHE:CD1	1:A:399:ILE:HG22	2.41	0.55
3:F:53:VAL:HG12	3:F:56:ASN:HB2	1.89	0.55
2:H:87:PHE:HB2	2:H:101:GLN:O	2.07	0.55
1:J:201:PHE:HE2	1:J:431:SER:HB2	1.72	0.55
1:D:261:LYS:O	1:D:261:LYS:NZ	2.31	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:146:SER:HA	3:C:149:LEU:HD21	1.89	0.54
3:F:119:PRO:HD2	3:F:121:PRO:HD3	1.90	0.54
1:G:225:LYS:HE3	2:H:50:ASP:HA	1.89	0.54
3:I:57:VAL:HG21	3:I:59:TRP:HH2	1.38	0.54
1:J:295:GLN:HA	1:J:298:LYS:HD2	1.88	0.54
3:F:149:LEU:HD22	3:F:207:LYS:HZ3	1.73	0.54
1:J:247:VAL:HG12	1:J:251:ILE:HD11	1.89	0.54
1:A:295:GLN:O	1:A:299:ILE:HG12	2.07	0.54
3:I:52:ASN:ND2	3:I:54:GLY:H	2.06	0.54
1:J:214:GLU:CG	1:J:311:LYS:H	2.19	0.54
1:A:358:ILE:HA	1:A:361:ILE:HG22	1.89	0.54
2:B:218:THR:HG22	2:B:219:CYS:O	2.08	0.54
3:F:140:SER:HB2	3:F:159:PHE:HB2	1.89	0.54
1:J:365:ILE:HD11	1:J:371:ILE:HD11	1.90	0.54
1:A:388:PHE:HE2	1:A:444:ASN:HB3	1.72	0.54
3:C:149:LEU:HD23	3:C:149:LEU:H	1.72	0.54
1:D:214:GLU:HG3	1:D:310:LEU:HA	1.90	0.54
1:G:419:LEU:HD12	1:G:423:TYR:HE2	1.73	0.54
1:J:325:ILE:O	1:J:329:LEU:HG	2.07	0.54
1:D:178:PHE:N	1:D:191:GLN:OE1	2.41	0.54
1:G:313:SER:O	1:G:314:GLN:HB2	2.08	0.54
2:B:66:TRP:HZ2	3:C:118:TYR:HD1	1.55	0.53
3:C:43:VAL:O	3:C:98:THR:HA	2.09	0.53
1:J:332:PHE:HE2	1:J:402:VAL:HG21	1.73	0.53
1:A:195:HIS:CD2	1:A:237:ILE:HG13	2.43	0.53
1:A:285:LEU:O	1:A:289:ASN:ND2	2.38	0.53
1:D:318:THR:HG22	1:D:419:LEU:HD13	1.90	0.53
2:E:108:GLU:N	2:E:108:GLU:OE1	2.41	0.53
3:F:130:ILE:HB	3:F:190:GLN:HE22	1.74	0.53
1:D:212:HIS:HB2	1:D:215:ILE:HG22	1.90	0.53
1:J:301:ASN:OD1	1:J:304:ARG:NH1	2.42	0.53
1:J:320:VAL:O	1:J:324:MET:N	2.41	0.53
1:D:218:TYR:HA	1:D:221:ILE:HD12	1.90	0.53
1:D:432:ILE:O	1:D:436:GLY:N	2.39	0.53
2:H:38:ASN:HB3	2:H:101:GLN:HG3	1.91	0.53
2:H:70:VAL:HG23	2:H:77:THR:HG22	1.91	0.53
1:A:374:ILE:HG12	1:A:457:LEU:HD22	1.91	0.53
2:B:66:TRP:NE1	3:C:120:TYR:H	2.07	0.53
1:D:318:THR:HB	1:D:321:TYR:HB3	1.91	0.53
1:D:340:TYR:HD2	1:D:388:PHE:HZ	1.57	0.52
2:B:143:PRO:HB3	2:B:169:TYR:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:PHE:HD1	1:A:399:ILE:HG22	1.73	0.52
1:G:322:PHE:O	1:G:326:LYS:HG2	2.09	0.52
1:G:356:ASP:OD1	1:G:356:ASP:N	2.39	0.52
2:H:87:PHE:CB	2:H:102:MET:HA	2.40	0.52
3:C:218:CYS:H	3:C:229:ILE:C	2.17	0.52
3:I:59:TRP:NE1	3:I:112:CYS:SG	2.83	0.52
1:J:226:VAL:HA	1:J:229:LEU:CD2	2.39	0.52
1:G:295:GLN:O	1:G:299:ILE:CD1	2.58	0.52
2:H:25:GLU:OE1	2:H:25:GLU:N	2.32	0.52
2:E:23:LEU:HD22	2:E:128:GLY:HA3	1.92	0.52
3:C:119:PRO:O	3:C:121:PRO:HD3	2.09	0.51
2:E:50:ASP:OD1	2:E:50:ASP:N	2.43	0.51
1:J:410:LYS:N	1:J:411:PRO:HD3	2.24	0.51
1:A:383:TYR:HD1	1:A:384:ILE:HD13	1.75	0.51
1:G:382:GLY:HA2	1:G:385:ILE:HD12	1.92	0.51
2:B:59:VAL:HB	2:B:62:LYS:HG2	1.91	0.51
3:C:173:LYS:HA	3:C:178:GLU:HA	1.91	0.51
2:E:86:ARG:NE	2:E:104:SER:O	2.42	0.51
1:G:415:THR:O	1:G:419:LEU:HD23	2.10	0.51
1:J:226:VAL:HA	1:J:229:LEU:HD23	1.92	0.51
1:J:416:LYS:O	1:J:420:LYS:HG3	2.10	0.51
1:G:261:LYS:HA	1:G:264:ARG:NH1	2.25	0.51
1:G:304:ARG:NH2	1:G:421:GLU:OE2	2.44	0.51
3:C:169:ASN:OD1	3:C:170:VAL:N	2.44	0.51
2:B:23:LEU:HD23	2:B:115:CYS:SG	2.50	0.51
3:C:113:GLN:HG2	3:C:114:GLN:N	2.24	0.51
1:G:177:TYR:OH	1:G:380:GLU:OE1	2.24	0.51
1:G:260:PHE:HE2	1:G:270:LYS:HD2	1.76	0.51
1:J:285:LEU:O	1:J:289:ASN:HB2	2.11	0.51
2:H:124:PHE:HE1	3:I:113:GLN:HE22	1.59	0.51
1:D:414:VAL:HG12	1:D:415:THR:H	1.76	0.51
3:F:183:VAL:HB	3:F:203:LEU:HD13	1.92	0.51
2:H:62:LYS:HD2	3:I:111:PHE:HB3	1.93	0.51
1:J:173:ASP:O	1:J:248:LYS:NZ	2.44	0.51
2:B:28:GLY:H	2:B:37:MET:HE3	1.75	0.50
3:F:205:LEU:HG	3:F:209:GLU:HG2	1.94	0.50
3:F:134:ASP:OD2	3:F:223:LYS:NZ	2.36	0.50
3:I:47:CYS:HB3	3:I:59:TRP:HE1	1.75	0.50
1:A:187:TYR:OH	1:A:297:LYS:HG2	2.10	0.50
1:A:308:GLU:HA	1:A:416:LYS:HE2	1.93	0.50
2:B:161:THR:HG22	2:B:206:THR:HG22	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:328:TYR:HB3	1:D:332:PHE:HE1	1.76	0.50
1:G:351:TYR:OH	1:G:378:GLN:O	2.27	0.50
1:J:200:TYR:CE1	1:J:434:LYS:HG3	2.46	0.50
1:A:267:TYR:CE2	1:A:372:ASN:HB3	2.47	0.50
1:A:314:GLN:HA	1:A:415:THR:HA	1.93	0.50
2:B:53:MET:HB3	2:B:98:LEU:HD22	1.93	0.50
3:F:173:LYS:HE2	3:F:219:GLU:HB2	1.93	0.50
2:B:193:VAL:O	2:B:200:THR:N	2.45	0.50
1:D:254:LEU:HD12	1:D:376:PHE:CD2	2.47	0.50
3:F:145:SER:O	3:F:149:LEU:HG	2.12	0.50
1:J:351:TYR:HB2	1:J:381:MET:HE2	1.94	0.50
1:A:380:GLU:O	1:A:384:ILE:HG12	2.12	0.50
2:B:32:GLN:HG3	2:B:33:PRO:HD2	1.93	0.50
2:H:80:LEU:O	2:H:84:LYS:N	2.45	0.50
2:H:84:LYS:HZ1	2:H:85:SER:H	1.59	0.50
3:I:59:TRP:HZ2	3:I:95:PHE:CB	2.25	0.50
2:B:110:THR:HG22	2:B:135:VAL:N	2.26	0.50
2:B:176:LEU:HD13	2:B:203:SER:HB3	1.94	0.50
3:C:49:ALA:O	3:C:93:THR:HG23	2.12	0.50
2:E:62:LYS:HA	2:E:63:GLY:C	2.36	0.50
2:E:107:SER:O	2:E:110:THR:OG1	2.30	0.50
3:F:216:TYR:HB2	3:F:231:LYS:HB3	1.93	0.50
3:I:47:CYS:O	3:I:95:PHE:HB2	2.12	0.50
1:D:267:TYR:CZ	1:D:372:ASN:HB3	2.47	0.50
2:B:165:LEU:HD13	2:B:202:SER:HB3	1.93	0.50
1:G:380:GLU:O	1:G:384:ILE:HG13	2.12	0.50
2:H:46:PHE:HB2	2:H:117:ARG:NH1	2.26	0.50
1:D:414:VAL:HG12	1:D:415:THR:N	2.27	0.49
1:G:419:LEU:HD23	1:G:419:LEU:H	1.76	0.49
1:A:279:GLU:O	1:A:283:GLY:N	2.42	0.49
2:B:195:GLN:O	2:B:197:ASP:N	2.46	0.49
3:C:208:ASP:OD1	3:C:209:GLU:HG2	2.12	0.49
1:G:208:TYR:CE2	1:G:426:ILE:HD11	2.47	0.49
1:G:253:ILE:HG22	1:G:270:LYS:HZ2	1.77	0.49
3:I:57:VAL:O	3:I:75:ALA:N	2.45	0.49
1:A:299:ILE:O	1:A:303:ILE:HG12	2.11	0.49
3:F:60:TYR:HD1	3:F:70:ALA:HA	1.76	0.49
1:J:222:HIS:HA	1:J:226:VAL:HB	1.93	0.49
1:A:354:GLY:O	1:A:358:ILE:HG13	2.12	0.49
1:D:263:GLN:OE1	1:D:265:GLU:N	2.35	0.49
2:E:180:SER:H	2:E:220:ASN:ND2	2.07	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:56:VAL:HG23	2:H:65:GLU:O	2.13	0.49
1:J:326:LYS:HD3	1:J:329:LEU:HD11	1.95	0.49
3:I:58:ALA:O	3:I:115:TYR:OH	2.16	0.49
1:J:358:ILE:HD11	1:J:371:ILE:HG23	1.94	0.49
1:A:196:PHE:HB2	1:A:199:ILE:HD13	1.94	0.49
2:B:36:SER:O	2:B:36:SER:OG	2.27	0.49
3:F:210:TYR:HE1	3:F:216:TYR:CE2	2.30	0.49
1:G:453:ILE:O	1:G:457:LEU:HD23	2.13	0.49
1:J:436:GLY:O	1:J:440:LEU:HG	2.12	0.49
2:E:65:GLU:N	3:F:122:PHE:HZ	2.09	0.48
2:B:55:TRP:CD1	2:B:100:LEU:HD23	2.48	0.48
2:B:171:PRO:O	2:B:223:HIS:NE2	2.46	0.48
1:D:206:LYS:HD3	1:D:222:HIS:CE1	2.48	0.48
2:E:53:MET:HB3	2:E:98:LEU:HD11	1.96	0.48
1:J:332:PHE:HA	1:J:433:PHE:CZ	2.48	0.48
1:D:446:ALA:O	1:D:450:LYS:HG2	2.13	0.48
1:A:236:ALA:HB1	1:A:288:LYS:HB3	1.94	0.48
3:C:37:THR:O	3:C:130:ILE:HA	2.12	0.48
3:C:132:ARG:HG3	3:C:133:ALA:N	2.29	0.48
2:H:43:ALA:HB3	2:H:48:PHE:CZ	2.48	0.48
1:A:269:VAL:O	1:A:273:GLN:HG3	2.14	0.48
1:A:274:TYR:OH	1:A:380:GLU:OE1	2.21	0.48
3:F:114:GLN:HG2	3:F:116:ASP:H	1.78	0.48
3:F:191:ASP:OD2	3:F:194:ASP:N	2.31	0.48
2:H:47:THR:O	2:H:117:ARG:NH2	2.47	0.48
1:J:303:ILE:HA	1:J:306:LEU:HD22	1.94	0.48
1:A:214:GLU:HG2	1:A:311:LYS:HG3	1.96	0.48
2:B:22:LYS:C	2:B:23:LEU:HD12	2.38	0.48
1:G:446:ALA:O	1:G:450:LYS:HG2	2.13	0.48
2:H:39:LEU:HB2	2:H:55:TRP:HZ3	1.78	0.48
1:J:313:SER:HB3	1:J:314:GLN:OE1	2.13	0.48
3:F:34:PHE:CD2	3:F:127:LYS:HB3	2.49	0.48
3:I:86:PHE:HZ	3:I:98:THR:O	1.96	0.48
1:J:181:SER:HA	1:J:188:LEU:HA	1.96	0.48
3:C:205:LEU:HD13	3:C:206:THR:O	2.14	0.48
1:D:448:ILE:O	1:D:452:LYS:HG2	2.14	0.48
2:E:57:ARG:NH2	2:E:109:ASP:OD1	2.47	0.48
3:I:52:ASN:ND2	3:I:54:GLY:N	2.62	0.47
3:I:78:ARG:NH2	3:I:84:ASP:OD1	2.44	0.47
1:J:394:LYS:HB3	1:J:432:ILE:HD12	1.95	0.47
3:C:61:GLN:HG3	3:C:110:TYR:HE1	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:69:SER:HB2	2:E:78:TYR:HB2	1.96	0.47
3:F:152:GLY:HA2	3:F:207:LYS:HG3	1.95	0.47
1:A:318:THR:N	1:A:412:LYS:O	2.41	0.47
1:A:381:MET:O	1:A:385:ILE:HG13	2.14	0.47
3:C:190:GLN:NE2	3:C:195:SER:HA	2.30	0.47
2:H:50:ASP:OD1	2:H:50:ASP:N	2.46	0.47
3:I:48:LYS:HG2	3:I:49:ALA:O	2.14	0.47
1:A:182:ASP:OD2	1:A:184:SER:OG	2.22	0.47
1:A:288:LYS:HD3	1:A:288:LYS:HA	1.71	0.47
1:D:224:THR:HG21	2:E:72:TYR:HD2	1.79	0.47
3:F:114:GLN:OE1	3:F:121:PRO:HD2	2.15	0.47
1:J:219:GLU:HB3	1:J:222:HIS:HB3	1.95	0.47
1:J:374:ILE:HB	1:J:457:LEU:HD22	1.96	0.47
3:C:174:ILE:N	3:C:177:SER:O	2.47	0.47
2:E:170:PHE:N	2:E:223:HIS:HE1	2.11	0.47
1:D:335:MET:HB2	1:D:335:MET:HE3	1.79	0.47
2:E:222:ALA:HB2	2:E:229:LYS:HD3	1.96	0.47
1:G:214:GLU:HB2	1:G:310:LEU:HG	1.96	0.47
1:G:312:CYS:SG	1:G:315:ASP:C	2.98	0.47
1:J:192:LEU:HD23	1:J:442:MET:HG2	1.96	0.47
1:J:351:TYR:HD2	1:J:352:LEU:HD22	1.79	0.47
1:A:344:ILE:HD13	1:A:392:LEU:HD23	1.97	0.47
3:F:157:VAL:HG13	3:F:202:THR:HG22	1.97	0.47
3:F:169:ASN:OD1	3:F:170:VAL:N	2.48	0.47
1:G:358:ILE:O	1:G:362:GLU:HB3	2.14	0.47
3:I:59:TRP:CD1	3:I:112:CYS:HB2	2.46	0.47
3:F:149:LEU:CD2	3:F:207:LYS:HZ3	2.28	0.47
1:G:397:HIS:CE1	1:G:401:GLN:HG3	2.50	0.47
1:J:175:ILE:HD11	1:J:244:LYS:HG3	1.96	0.47
1:G:386:ASP:O	1:G:390:TYR:N	2.45	0.46
2:H:115:CYS:O	2:H:128:GLY:N	2.48	0.46
1:A:170:ASN:OD1	1:A:170:ASN:N	2.48	0.46
2:B:222:ALA:O	2:B:224:PRO:HD3	2.16	0.46
3:C:59:TRP:CH2	3:C:112:CYS:HB3	2.50	0.46
1:G:240:CYS:O	1:G:244:LYS:N	2.38	0.46
1:G:299:ILE:O	1:G:303:ILE:HG13	2.14	0.46
1:J:419:LEU:HD13	1:J:419:LEU:O	2.15	0.46
2:B:111:ALA:HB3	2:B:113:TYR:CZ	2.50	0.46
1:D:312:CYS:O	1:D:416:LYS:HE2	2.15	0.46
3:F:112:CYS:O	3:F:122:PHE:HB2	2.16	0.46
2:H:127:TRP:CE3	3:I:68:PRO:HG2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:182:ASP:HB3	1:D:286:LEU:HD11	1.98	0.46
3:C:179:ARG:NE	3:C:181:ASN:O	2.49	0.46
1:G:251:ILE:HG22	1:G:255:GLU:OE2	2.16	0.46
1:G:247:VAL:HG21	1:G:281:PHE:HB2	1.98	0.46
2:H:61:ASP:HA	2:H:62:LYS:HA	1.56	0.46
3:I:53:VAL:HG21	3:I:116:ASP:OD2	2.06	0.46
3:C:52:ASN:OD1	3:C:92:GLY:HA2	2.16	0.46
3:C:57:VAL:HA	3:C:113:GLN:O	2.15	0.46
1:G:316:CYS:HB2	1:G:419:LEU:HD22	1.97	0.46
3:I:56:ASN:C	3:I:74:SER:HA	2.41	0.46
2:B:139:LYS:H	2:B:139:LYS:HD3	1.80	0.46
1:D:275:GLU:O	1:D:279:GLU:HG2	2.16	0.46
2:E:171:PRO:CD	2:E:225:ALA:HB3	2.46	0.46
1:J:214:GLU:OE2	1:J:311:LYS:HB3	2.16	0.46
2:B:35:SER:HB2	2:B:103:SER:HA	1.97	0.45
3:C:155:SER:HB3	3:C:204:THR:HG22	1.98	0.45
2:E:109:ASP:C	2:E:111:ALA:H	2.24	0.45
1:J:207:ARG:HD2	1:J:207:ARG:HA	1.80	0.45
1:J:258:GLN:O	1:J:262:THR:HG23	2.15	0.45
1:J:322:PHE:O	1:J:325:ILE:HG22	2.16	0.45
2:E:58:GLN:CD	2:E:62:LYS:HE3	2.41	0.45
1:G:256:ASN:HD21	1:G:259:LYS:HB2	1.81	0.45
1:G:393:GLN:O	1:G:393:GLN:NE2	2.48	0.45
1:G:400:ASP:OD1	1:G:401:GLN:N	2.50	0.45
1:J:225:LYS:C	1:J:229:LEU:HD23	2.40	0.45
2:B:41:CYS:O	2:B:97:ILE:HA	2.16	0.45
3:C:171:LYS:HE3	3:C:221:THR:OG1	2.16	0.45
2:E:62:LYS:HB2	3:F:111:PHE:CZ	2.52	0.45
3:F:165:PRO:HG3	3:F:223:LYS:HE3	1.99	0.45
1:G:322:PHE:HZ	1:G:411:PRO:HD2	1.81	0.45
2:H:48:PHE:HB3	2:H:91:ARG:NH2	2.31	0.45
1:J:212:HIS:CE1	1:J:216:LYS:HA	2.51	0.45
2:E:212:TRP:CZ2	2:E:236:PRO:HB3	2.52	0.45
3:F:190:GLN:HG3	3:F:197:TYR:CE1	2.52	0.45
1:G:170:ASN:OD1	1:G:171:THR:N	2.50	0.45
1:A:378:GLN:HA	1:A:381:MET:HE3	1.99	0.45
2:B:23:LEU:HA	2:B:42:THR:O	2.16	0.45
2:B:204:SER:HB3	3:C:159:PHE:CE2	2.52	0.45
1:G:451:GLU:O	1:G:454:VAL:HG12	2.17	0.45
2:H:55:TRP:HE1	2:H:98:LEU:HD23	1.82	0.45
1:J:197:SER:O	1:J:431:SER:OG	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:190:SER:HB2	1:G:435:PHE:HZ	1.82	0.45
1:G:442:MET:HE2	1:G:442:MET:HB2	1.86	0.45
3:I:45:VAL:O	3:I:97:LEU:HB3	2.16	0.45
1:J:347:TYR:OH	1:J:451:GLU:HB2	2.16	0.45
2:B:73:ASP:OD1	2:B:74:GLY:N	2.50	0.45
2:E:64:LEU:HD23	2:E:64:LEU:HA	1.82	0.45
1:J:258:GLN:HA	1:J:261:LYS:HG3	1.99	0.45
1:J:332:PHE:CD2	1:J:399:ILE:HG22	2.52	0.45
1:A:267:TYR:OH	1:A:376:PHE:HB2	2.16	0.45
1:A:307:LEU:HB2	1:A:420:LYS:HE2	1.98	0.45
1:G:312:CYS:SG	1:G:316:CYS:N	2.90	0.45
1:G:415:THR:O	1:G:418:ARG:N	2.48	0.45
2:H:46:PHE:HB2	2:H:117:ARG:HH12	1.82	0.45
2:E:172:GLU:N	2:E:173:PRO:HD2	2.32	0.44
1:G:220:ASP:OD2	2:H:71:SER:OG	2.34	0.44
2:B:21:VAL:HA	2:B:44:SER:O	2.17	0.44
3:F:154:ALA:O	3:F:205:LEU:N	2.31	0.44
1:G:308:GLU:HA	1:G:416:LYS:HZ3	1.82	0.44
1:G:362:GLU:OE2	1:G:362:GLU:N	2.50	0.44
3:C:28:MET:HE2	3:C:53:VAL:HG22	1.97	0.44
1:J:225:LYS:O	1:J:229:LEU:N	2.41	0.44
1:J:251:ILE:HG22	1:J:255:GLU:OE1	2.16	0.44
1:J:370:THR:HG23	1:J:371:ILE:HD12	2.00	0.44
3:C:52:ASN:OD1	3:C:54:GLY:N	2.37	0.44
3:F:44:SER:HA	3:F:97:LEU:O	2.18	0.44
1:A:293:LEU:O	1:A:297:LYS:HG3	2.18	0.44
1:A:328:TYR:O	1:A:332:PHE:N	2.43	0.44
1:D:336:PRO:HD2	1:D:340:TYR:HE1	1.81	0.44
3:F:38:SER:H	3:F:41:VAL:HG22	1.83	0.44
1:G:322:PHE:CZ	1:G:411:PRO:HD2	2.52	0.44
1:J:287:ASN:OD1	1:J:287:ASN:N	2.43	0.44
1:A:388:PHE:CE2	1:A:444:ASN:HB3	2.51	0.44
3:I:86:PHE:CZ	3:I:98:THR:C	2.96	0.44
1:J:297:LYS:HB2	1:J:297:LYS:HE3	1.79	0.44
1:J:394:LYS:HA	1:J:394:LYS:HD2	1.86	0.44
2:B:68:ALA:HB2	2:B:79:TYR:HD1	1.83	0.44
1:D:469:ILE:O	1:D:470:SER:HB3	2.17	0.44
2:E:171:PRO:HB2	2:E:173:PRO:HD2	2.00	0.44
1:G:417:ASN:O	1:G:421:GLU:HG2	2.17	0.44
1:J:363:LYS:H	1:J:363:LYS:HG3	1.67	0.44
1:J:440:LEU:HD12	1:J:441:ASN:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:ARG:HA	1:A:264:ARG:NE	2.32	0.44
1:A:266:SER:HB3	1:A:269:VAL:HG22	2.00	0.44
3:I:61:GLN:HG3	3:I:62:GLN:O	2.18	0.44
2:B:59:VAL:O	2:B:62:LYS:HB2	2.18	0.44
3:C:147:GLU:O	3:C:150:THR:OG1	2.33	0.44
1:D:359:ARG:HA	1:D:362:GLU:HG3	2.00	0.44
1:G:331:ASP:OD2	1:G:430:TYR:OH	2.35	0.44
1:J:220:ASP:OD1	1:J:224:THR:OG1	2.36	0.44
2:B:59:VAL:HG12	2:B:60:PRO:HD2	1.99	0.43
2:B:71:SER:C	2:B:91:ARG:NH1	2.76	0.43
2:E:58:GLN:C	2:E:111:ALA:HB1	2.42	0.43
3:I:52:ASN:ND2	3:I:54:GLY:CA	2.81	0.43
1:J:178:PHE:CE1	1:J:278:LYS:HE3	2.53	0.43
1:J:259:LYS:O	1:J:262:THR:OG1	2.26	0.43
1:J:307:LEU:HD13	1:J:420:LYS:HG2	1.99	0.43
1:A:244:LYS:HB2	1:A:281:PHE:CE1	2.53	0.43
1:A:390:TYR:HD1	1:A:391:HIS:HD2	1.66	0.43
2:B:113:TYR:O	2:B:130:GLY:HA2	2.18	0.43
2:B:163:GLY:HA2	2:B:204:SER:HA	1.99	0.43
2:B:198:LEU:HD23	2:B:198:LEU:HA	1.79	0.43
3:C:28:MET:HE3	3:C:49:ALA:HB3	1.98	0.43
3:C:74:SER:O	3:C:76:SER:N	2.45	0.43
2:H:39:LEU:HB2	2:H:55:TRP:CZ3	2.52	0.43
1:J:371:ILE:HD12	1:J:371:ILE:N	2.33	0.43
2:B:55:TRP:CE3	2:B:113:TYR:HB3	2.54	0.43
2:B:111:ALA:HB3	2:B:113:TYR:CE2	2.54	0.43
2:B:173:PRO:O	2:B:224:PRO:HD2	2.18	0.43
2:E:62:LYS:H	2:E:62:LYS:HD2	1.83	0.43
3:F:59:TRP:HB2	3:F:72:ILE:HB	1.98	0.43
1:G:310:LEU:HD22	1:G:311:LYS:HG2	2.01	0.43
3:I:61:GLN:HG3	3:I:62:GLN:N	2.33	0.43
1:A:365:ILE:H	1:A:365:ILE:HG13	1.67	0.43
2:B:141:THR:N	2:B:170:PHE:O	2.45	0.43
1:D:354:GLY:O	1:D:358:ILE:HG13	2.18	0.43
3:F:147:GLU:O	3:F:150:THR:HG22	2.18	0.43
2:B:25:GLU:CD	2:B:130:GLY:H	2.27	0.43
2:B:218:THR:HG23	2:B:232:LYS:C	2.43	0.43
3:C:48:LYS:O	3:C:49:ALA:HB2	2.17	0.43
1:D:347:TYR:CE2	1:D:447:LEU:HD22	2.54	0.43
1:G:344:ILE:HD12	1:G:344:ILE:HA	1.89	0.43
1:G:364:GLN:N	1:G:364:GLN:OE1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:48:LYS:HG2	3:I:49:ALA:N	2.33	0.43
1:A:387:ARG:HA	1:A:387:ARG:HD3	1.68	0.43
1:D:206:LYS:HD3	1:D:222:HIS:ND1	2.33	0.43
1:D:374:ILE:O	1:D:378:GLN:HG3	2.18	0.43
2:E:188:HIS:NE2	3:F:188:THR:HG21	2.34	0.43
2:E:212:TRP:CH2	2:E:236:PRO:HB3	2.54	0.43
1:G:172:THR:HG22	1:G:449:HIS:CE1	2.53	0.43
1:G:247:VAL:O	1:G:251:ILE:HG13	2.19	0.43
1:J:410:LYS:N	1:J:411:PRO:CD	2.82	0.43
2:B:24:VAL:O	2:B:41:CYS:HA	2.19	0.43
1:D:419:LEU:HA	1:D:422:TYR:HB2	2.01	0.43
1:G:325:ILE:HG22	1:G:426:ILE:HG21	2.00	0.43
3:C:39:ILE:HA	3:C:40:GLY:HA2	1.54	0.43
3:C:179:ARG:HD3	3:C:203:LEU:HD11	1.99	0.43
1:D:441:ASN:O	1:D:445:LYS:HG3	2.19	0.43
1:G:422:TYR:HA	1:G:425:ASN:ND2	2.34	0.43
2:H:54:ALA:O	2:H:116:THR:OG1	2.21	0.43
3:C:175:ASP:HA	3:C:213:HIS:ND1	2.34	0.43
1:D:176:ASP:HB2	1:D:191:GLN:HG2	2.00	0.43
1:G:273:GLN:O	1:G:277:LYS:HG2	2.19	0.43
1:J:336:PRO:HB2	1:J:339:ASN:HB2	2.01	0.43
2:E:86:ARG:HG2	2:E:104:SER:HB2	2.00	0.43
1:G:266:SER:HB3	1:G:269:VAL:HG22	2.00	0.43
2:B:38:ASN:OD1	2:B:101:GLN:HA	2.19	0.42
1:G:260:PHE:CE2	1:G:270:LYS:HD2	2.53	0.42
1:G:328:TYR:HE2	1:G:426:ILE:HD13	1.84	0.42
1:J:279:GLU:O	1:J:283:GLY:N	2.45	0.42
1:J:446:ALA:O	1:J:450:LYS:HG2	2.19	0.42
2:B:59:VAL:H	2:B:62:LYS:CG	2.32	0.42
2:E:174:VAL:HG13	2:E:222:ALA:O	2.18	0.42
1:A:275:GLU:O	1:A:279:GLU:HG2	2.19	0.42
2:B:58:GLN:N	2:B:112:THR:O	2.52	0.42
1:D:222:HIS:HA	1:D:226:VAL:CG1	2.48	0.42
1:G:373:ALA:O	1:G:377:THR:HG23	2.18	0.42
2:H:40:SER:HA	2:H:98:LEU:O	2.18	0.42
2:H:55:TRP:CD1	2:H:115:CYS:HA	2.54	0.42
1:D:227:ASN:OD1	1:D:228:SER:N	2.52	0.42
1:G:260:PHE:HE1	1:G:264:ARG:HA	1.83	0.42
1:A:316:CYS:HB3	1:A:419:LEU:HD22	2.01	0.42
1:A:374:ILE:CG1	1:A:457:LEU:HD22	2.49	0.42
1:D:336:PRO:HD2	1:D:340:TYR:CE1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:210:SER:O	2:E:213:PRO:HD2	2.19	0.42
1:J:242:ARG:O	1:J:246:THR:HG23	2.19	0.42
1:J:275:GLU:O	1:J:279:GLU:HG2	2.19	0.42
2:B:62:LYS:HA	2:B:62:LYS:HD3	1.68	0.42
1:D:183:GLU:OE1	1:D:183:GLU:N	2.46	0.42
1:D:194:PRO:C	1:D:196:PHE:H	2.26	0.42
1:D:254:LEU:HD12	1:D:376:PHE:CG	2.55	0.42
2:H:57:ARG:HB3	2:H:113:TYR:CD1	2.54	0.42
3:I:62:GLN:HB3	3:I:68:PRO:HA	2.01	0.42
1:J:405:LEU:O	1:J:409:VAL:HG22	2.19	0.42
1:A:368:PRO:HB2	1:A:372:ASN:OD1	2.20	0.42
2:B:59:VAL:CG1	2:B:60:PRO:HD2	2.49	0.42
3:C:190:GLN:HE21	3:C:195:SER:HA	1.84	0.42
3:F:57:VAL:HA	3:F:113:GLN:O	2.20	0.42
1:J:332:PHE:CE2	1:J:402:VAL:HG21	2.51	0.42
1:J:374:ILE:CA	1:J:457:LEU:HD13	2.50	0.42
1:A:451:GLU:O	1:A:454:VAL:HG12	2.20	0.42
2:B:27:GLU:HA	2:B:28:GLY:HA3	1.76	0.42
2:B:180:SER:H	2:B:220:ASN:ND2	2.18	0.42
1:G:232:GLU:CG	1:G:292:ASN:HD21	2.33	0.42
2:H:83:LEU:HD12	2:H:87:PHE:CZ	2.55	0.42
2:H:98:LEU:HD12	2:H:99:ASN:H	1.83	0.42
1:J:460:GLU:O	1:J:464:HIS:N	2.52	0.42
1:A:310:LEU:O	1:A:416:LYS:NZ	2.34	0.42
1:A:357:MET:HE2	1:A:358:ILE:HG12	2.02	0.42
3:C:28:MET:HB2	3:C:50:SER:N	2.16	0.42
3:C:226:THR:C	3:C:228:PRO:HD3	2.45	0.42
2:E:66:TRP:HZ3	3:F:118:TYR:CE2	2.38	0.42
1:J:192:LEU:CD2	1:J:442:MET:HG2	2.50	0.42
1:J:340:TYR:CZ	1:J:440:LEU:HD13	2.55	0.42
1:A:340:TYR:HD1	1:A:388:PHE:HZ	1.67	0.42
1:A:393:GLN:O	1:A:397:HIS:N	2.47	0.42
2:E:125:ASP:C	2:E:126:TYR:HD1	2.28	0.42
1:G:214:GLU:OE2	1:G:309:LYS:HE3	2.20	0.42
1:G:335:MET:HE2	1:G:335:MET:HB3	1.93	0.42
1:G:449:HIS:O	1:G:453:ILE:HG22	2.20	0.42
3:C:113:GLN:HE21	3:C:120:TYR:HB3	1.85	0.41
1:G:468:ARG:NE	1:G:468:ARG:HA	2.34	0.41
2:H:83:LEU:HD12	2:H:87:PHE:CE2	2.54	0.41
1:J:289:ASN:OD1	1:J:292:ASN:ND2	2.47	0.41
1:D:309:LYS:NZ	2:E:122:ASP:OD2	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:260:PHE:O	1:G:264:ARG:HG2	2.20	0.41
1:J:286:LEU:O	1:J:290:ARG:HB2	2.20	0.41
1:J:318:THR:OG1	1:J:412:LYS:HA	2.21	0.41
2:B:118:VAL:HG22	2:B:124:PHE:CD1	2.56	0.41
1:D:244:LYS:HB2	1:D:281:PHE:CE1	2.55	0.41
3:F:148:GLN:OE1	3:F:155:SER:OG	2.28	0.41
1:G:232:GLU:HG2	1:G:292:ASN:HD21	1.84	0.41
3:I:53:VAL:HG23	3:I:54:GLY:N	2.35	0.41
2:B:227:SER:O	2:B:227:SER:OG	2.29	0.41
3:C:42:ARG:HA	3:C:99:ILE:O	2.20	0.41
1:D:243:ALA:O	1:D:247:VAL:HG23	2.21	0.41
1:D:264:ARG:HD2	1:D:265:GLU:N	2.35	0.41
1:G:253:ILE:HG22	1:G:270:LYS:NZ	2.35	0.41
1:J:419:LEU:HD22	1:J:422:TYR:HB2	2.02	0.41
2:E:148:LEU:HD11	3:F:142:PHE:HB3	2.03	0.41
3:F:164:TYR:CD1	3:F:165:PRO:HA	2.55	0.41
1:G:397:HIS:O	1:G:397:HIS:ND1	2.54	0.41
2:H:90:SER:N	2:H:99:ASN:O	2.30	0.41
1:A:338:GLU:OE1	1:A:338:GLU:N	2.31	0.41
2:E:169:TYR:OH	2:E:201:LEU:HD23	2.19	0.41
2:E:223:HIS:HB3	2:E:226:SER:HB2	2.02	0.41
1:J:176:ASP:O	1:J:442:MET:HE1	2.20	0.41
2:B:180:SER:H	2:B:220:ASN:HD21	1.69	0.41
1:D:340:TYR:HD2	1:D:388:PHE:CZ	2.37	0.41
1:D:394:LYS:HA	1:D:394:LYS:HD3	1.77	0.41
2:E:149:ALA:O	3:F:143:PRO:HD2	2.21	0.41
1:G:337:TYR:CE2	1:G:396:LYS:HD3	2.56	0.41
1:G:366:ASP:HB3	1:G:468:ARG:CD	2.50	0.41
3:I:55:THR:HA	3:I:95:PHE:CE1	2.55	0.41
3:I:112:CYS:SG	3:I:112:CYS:O	2.78	0.41
1:A:177:TYR:HA	1:A:442:MET:HE1	2.03	0.41
1:A:329:LEU:HA	1:A:332:PHE:HB2	2.03	0.41
2:B:35:SER:OG	2:B:36:SER:N	2.53	0.41
1:D:295:GLN:O	1:D:299:ILE:HG13	2.20	0.41
1:G:340:TYR:CE1	1:G:440:LEU:HB3	2.55	0.41
2:H:79:TYR:CE1	2:H:89:ILE:HG22	2.56	0.41
1:A:399:ILE:O	1:A:403:THR:OG1	2.19	0.41
1:A:441:ASN:O	1:A:444:ASN:ND2	2.54	0.41
2:B:139:LYS:HD3	2:B:139:LYS:N	2.36	0.41
2:B:179:ASN:CA	2:B:217:ILE:HG21	2.40	0.41
2:B:218:THR:HG23	2:B:232:LYS:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:35:SER:HB3	2:E:105:LEU:HD11	2.03	0.41
2:E:58:GLN:NE2	2:E:62:LYS:HE3	2.36	0.41
2:E:143:PRO:HD3	2:E:223:HIS:ND1	2.36	0.41
3:F:219:GLU:HG3	3:F:228:PRO:HB3	2.01	0.41
1:G:197:SER:HB2	1:G:435:PHE:CD2	2.56	0.41
1:G:298:LYS:O	1:G:302:GLU:HG3	2.20	0.41
1:G:368:PRO:HA	1:G:371:ILE:HG12	2.02	0.41
1:G:386:ASP:OD1	1:G:386:ASP:N	2.45	0.41
1:G:394:LYS:HA	1:G:394:LYS:HD3	1.93	0.41
2:H:84:LYS:C	2:H:86:ARG:H	2.29	0.41
1:A:309:LYS:HZ1	2:B:122:ASP:CG	2.26	0.41
1:D:217:ARG:HH21	1:D:221:ILE:HD11	1.86	0.41
1:D:347:TYR:OH	1:D:451:GLU:HB2	2.21	0.41
3:F:54:GLY:H	3:F:92:GLY:HA2	1.86	0.41
1:J:217:ARG:HE	1:J:306:LEU:HG	1.83	0.41
1:J:254:LEU:HB3	1:J:376:PHE:CE2	2.56	0.41
1:J:427:GLY:O	1:J:430:TYR:N	2.54	0.41
1:A:441:ASN:HA	1:A:444:ASN:HD21	1.86	0.40
3:C:55:THR:O	3:C:74:SER:HA	2.21	0.40
1:D:180:ILE:HD12	1:D:282:ARG:HB3	2.03	0.40
1:D:253:ILE:HG23	1:D:260:PHE:HA	2.02	0.40
1:G:306:LEU:HG	2:H:120:VAL:HG12	2.02	0.40
1:G:385:ILE:O	1:G:389:GLU:N	2.40	0.40
1:G:410:LYS:HA	1:G:410:LYS:HD2	1.91	0.40
1:J:366:ASP:OD1	1:J:367:ASN:N	2.54	0.40
2:B:174:VAL:HG13	2:B:222:ALA:O	2.21	0.40
1:D:215:ILE:HG13	1:D:306:LEU:HG	2.03	0.40
1:G:359:ARG:CD	1:J:290:ARG:HE	2.35	0.40
1:J:254:LEU:HD11	1:J:270:LYS:CB	2.50	0.40
1:J:299:ILE:O	1:J:303:ILE:HG13	2.21	0.40
1:D:385:ILE:O	1:D:389:GLU:HG3	2.21	0.40
1:A:447:LEU:HD23	1:A:447:LEU:HA	1.83	0.40
1:G:269:VAL:O	1:G:273:GLN:HG2	2.21	0.40
3:I:55:THR:HA	3:I:95:PHE:HE1	1.85	0.40
1:A:454:VAL:O	1:A:458:LEU:HG	2.22	0.40
2:B:98:LEU:HD23	2:B:115:CYS:SG	2.61	0.40
2:B:149:ALA:O	3:C:143:PRO:HD2	2.22	0.40
3:C:61:GLN:CB	3:C:71:LEU:HD11	2.51	0.40
2:H:81:ASP:HA	2:H:84:LYS:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	297/307 (97%)	289 (97%)	8 (3%)	0	100	100
1	D	300/307 (98%)	295 (98%)	4 (1%)	1 (0%)	36	65
1	G	300/307 (98%)	289 (96%)	11 (4%)	0	100	100
1	J	295/307 (96%)	284 (96%)	10 (3%)	1 (0%)	36	65
2	B	205/254 (81%)	193 (94%)	10 (5%)	2 (1%)	12	41
2	E	204/254 (80%)	192 (94%)	10 (5%)	2 (1%)	12	41
2	H	111/254 (44%)	107 (96%)	4 (4%)	0	100	100
3	C	195/238 (82%)	183 (94%)	10 (5%)	2 (1%)	12	41
3	F	192/238 (81%)	180 (94%)	12 (6%)	0	100	100
3	I	61/238 (26%)	56 (92%)	5 (8%)	0	100	100
All	All	2160/2704 (80%)	2068 (96%)	84 (4%)	8 (0%)	30	59

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	179	ASN
3	C	52	ASN
3	C	228	PRO
1	D	336	PRO
2	E	224	PRO
2	B	64	LEU
2	B	196	SER
1	J	410	LYS

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	281/286 (98%)	281 (100%)	0	100	100
1	D	284/286 (99%)	284 (100%)	0	100	100
1	G	284/286 (99%)	284 (100%)	0	100	100
1	J	279/286 (98%)	279 (100%)	0	100	100
2	B	187/227 (82%)	187 (100%)	0	100	100
2	E	186/227 (82%)	186 (100%)	0	100	100
2	H	101/227 (44%)	101 (100%)	0	100	100
3	C	175/211 (83%)	175 (100%)	0	100	100
3	F	175/211 (83%)	175 (100%)	0	100	100
3	I	58/211 (28%)	56 (97%)	2 (3%)	32	59
All	All	2010/2458 (82%)	2008 (100%)	2 (0%)	88	89

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

Mol	Chain	Res	Type
3	I	51	GLN
3	I	112	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	372	ASN
1	A	401	GLN
1	A	444	ASN
3	C	32	GLN
3	C	62	GLN
1	D	289	ASN
1	D	317	GLN
1	D	364	GLN
1	D	378	GLN
1	D	425	ASN
1	D	444	ASN
2	E	32	GLN
2	E	38	ASN
2	E	58	GLN
2	E	179	ASN

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Mol	Chain	Res	Type
3	F	190	GLN
1	G	227	ASN
1	G	287	ASN
1	G	292	ASN
1	G	372	ASN
1	G	413	GLN
1	G	428	ASN
1	G	455	HIS
3	I	113	GLN
1	J	339	ASN
1	J	441	ASN
1	J	455	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	299/307 (97%)	0.63	31 (10%)	11 10	28, 67, 115, 149	0
1	D	302/307 (98%)	0.72	30 (9%)	13 11	36, 72, 116, 173	0
1	G	302/307 (98%)	0.95	40 (13%)	7 7	41, 85, 140, 181	0
1	J	297/307 (96%)	1.09	49 (16%)	4 5	69, 109, 149, 189	0
2	B	209/254 (82%)	0.81	23 (11%)	10 10	31, 72, 135, 165	0
2	E	208/254 (81%)	0.95	31 (14%)	5 6	34, 86, 130, 144	0
2	H	113/254 (44%)	0.72	14 (12%)	8 7	56, 86, 153, 180	0
3	C	199/238 (83%)	1.00	30 (15%)	5 5	34, 79, 130, 176	0
3	F	198/238 (83%)	0.85	23 (11%)	9 9	38, 74, 130, 170	0
3	I	67/238 (28%)	1.12	10 (14%)	5 6	81, 117, 163, 192	0
All	All	2194/2704 (81%)	0.87	281 (12%)	7 7	28, 82, 140, 192	0

All (281) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	187	TYR	7.7
3	F	37	THR	5.9
1	G	208	TYR	5.8
2	E	164	CYS	5.8
3	F	108	ALA	5.6
1	G	279	GLU	5.5
1	J	303	ILE	4.9
1	A	215	ILE	4.8
1	D	371	ILE	4.8
2	B	63	GLY	4.7
1	A	457	LEU	4.6
1	G	408	GLY	4.6
3	C	28	MET	4.5

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Mol	Chain	Res	Type	RSRZ
3	F	27	VAL	4.5
1	G	297	LYS	4.3
1	J	187	TYR	4.3
1	D	389	GLU	4.3
3	I	59	TRP	4.2
2	E	236	PRO	4.2
2	E	214	SER	4.1
3	C	154	ALA	4.1
2	B	65	GLU	4.1
1	D	205	PHE	4.0
1	J	217	ARG	4.0
1	A	453	ILE	3.9
1	J	213	THR	3.9
3	F	29	THR	3.9
2	B	146	TYR	3.8
3	C	143	PRO	3.8
2	E	64	LEU	3.8
2	B	95	LYS	3.8
3	C	39	ILE	3.8
1	J	207	ARG	3.8
3	I	71	LEU	3.7
1	J	424	PHE	3.7
1	D	451	GLU	3.7
1	J	211	TYR	3.7
1	D	346	GLN	3.7
2	E	162	LEU	3.7
3	C	31	SER	3.7
2	H	58	GLN	3.7
2	H	48	PHE	3.7
3	C	124	GLY	3.7
1	D	173	ASP	3.6
1	D	464	HIS	3.6
3	C	217	THR	3.6
2	E	187	VAL	3.5
1	G	215	ILE	3.4
2	H	67	VAL	3.4
2	E	37	MET	3.4
1	J	210	SER	3.4
1	J	276	GLU	3.3
3	C	83	PRO	3.2
3	C	149	LEU	3.2
3	I	87	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	G	267	TYR	3.2
1	J	181	SER	3.2
3	F	167	ASP	3.2
2	B	136	SER	3.2
1	G	283	GLY	3.1
1	A	228	SER	3.1
3	I	60	TYR	3.1
2	E	35	SER	3.1
3	C	38	SER	3.1
1	J	357	MET	3.1
1	J	428	ASN	3.0
1	G	321	TYR	3.0
3	C	110	TYR	3.0
1	A	464	HIS	3.0
2	B	147	PRO	3.0
1	G	448	ILE	3.0
3	F	110	TYR	3.0
3	C	167	ASP	3.0
1	J	283	GLY	3.0
3	F	61	GLN	3.0
1	G	198	ASN	3.0
1	G	205	PHE	3.0
3	C	34	PHE	3.0
1	A	187	TYR	3.0
1	D	369	VAL	3.0
1	G	409	VAL	2.9
3	C	123	GLY	2.9
2	B	64	LEU	2.9
1	G	369	VAL	2.9
1	J	409	VAL	2.9
1	D	462	PHE	2.9
1	D	318	THR	2.9
1	J	356	ASP	2.8
3	I	49	ALA	2.8
1	D	358	ILE	2.8
1	A	173	ASP	2.8
2	B	62	LYS	2.8
1	J	199	ILE	2.8
3	F	31	SER	2.8
1	A	203	ASP	2.8
2	B	161	THR	2.8
1	A	322	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	G	196	PHE	2.8
1	D	254	LEU	2.8
1	G	307	LEU	2.8
1	J	221	ILE	2.8
1	G	185	ASN	2.8
1	G	293	LEU	2.8
2	H	136	SER	2.7
3	F	128	LEU	2.7
3	I	119	PRO	2.7
1	A	460	GLU	2.7
2	E	34	GLY	2.7
1	G	463	GLY	2.7
1	A	270	LYS	2.7
1	J	297	LYS	2.7
1	D	303	ILE	2.7
1	A	205	PHE	2.7
1	J	229	LEU	2.7
1	J	306	LEU	2.7
1	J	405	LEU	2.7
1	J	414	VAL	2.7
1	G	194	PRO	2.6
1	J	250	LEU	2.6
1	J	426	ILE	2.6
2	B	125	ASP	2.6
1	A	324	MET	2.6
3	C	32	GLN	2.6
3	F	218	CYS	2.6
3	F	157	VAL	2.6
1	J	365	ILE	2.6
3	C	223	LYS	2.6
2	E	161	THR	2.6
3	C	46	THR	2.6
3	C	218	CYS	2.6
3	C	115	TYR	2.6
3	C	30	GLN	2.6
1	A	441	ASN	2.6
1	G	465	LEU	2.6
1	J	358	ILE	2.6
2	H	36	SER	2.6
2	B	94	VAL	2.6
1	J	423	TYR	2.5
1	D	190	SER	2.5

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Mol	Chain	Res	Type	RSRZ
3	F	225	SER	2.5
2	E	163	GLY	2.5
1	D	188	LEU	2.5
1	J	182	ASP	2.5
1	D	307	LEU	2.5
2	E	104	SER	2.5
2	H	72	TYR	2.5
2	E	160	VAL	2.5
1	G	310	LEU	2.5
1	A	208	TYR	2.5
1	D	222	HIS	2.5
2	H	52	TYR	2.5
1	A	346	GLN	2.5
1	J	439	SER	2.4
2	B	123	SER	2.4
2	H	55	TRP	2.4
1	A	371	ILE	2.4
1	J	432	ILE	2.4
1	J	275	GLU	2.4
1	J	216	LYS	2.4
3	F	160	LEU	2.4
1	A	366	ASP	2.4
1	G	220	ASP	2.4
3	C	133	ALA	2.4
3	F	109	GLU	2.4
2	E	85	SER	2.4
3	C	100	SER	2.4
2	B	19	CYS	2.4
1	G	257	PRO	2.4
2	B	228	THR	2.4
2	H	33	PRO	2.4
1	J	285	LEU	2.4
2	B	162	LEU	2.4
1	D	347	TYR	2.4
2	E	122	ASP	2.4
2	E	176	LEU	2.4
2	E	206	THR	2.3
2	E	211	THR	2.3
2	B	149	ALA	2.3
3	F	229	ILE	2.3
2	E	125	ASP	2.3
1	J	183	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	253	ILE	2.3
1	D	266	SER	2.3
3	C	220	ALA	2.3
2	E	199	TYR	2.3
2	H	78	TYR	2.3
2	H	64	LEU	2.3
3	C	214	ASN	2.3
1	A	179	ASP	2.3
3	C	29	THR	2.3
2	B	55	TRP	2.3
1	G	457	LEU	2.3
1	G	186	TYR	2.3
1	J	196	PHE	2.3
2	B	216	SER	2.3
1	J	274	TYR	2.3
3	C	122	PHE	2.3
3	I	111	PHE	2.3
2	B	115	CYS	2.3
3	C	171	LYS	2.3
1	J	220	ASP	2.3
1	D	306	LEU	2.2
1	D	458	LEU	2.2
1	A	267	TYR	2.2
1	D	211	TYR	2.2
2	E	62	LYS	2.2
3	C	186	SER	2.2
1	A	264	ARG	2.2
2	H	38	ASN	2.2
3	C	113	GLN	2.2
2	E	115	CYS	2.2
1	J	179	ASP	2.2
2	B	61	ASP	2.2
2	E	150	PRO	2.2
2	B	201	LEU	2.2
3	F	125	GLY	2.2
1	J	369	VAL	2.2
1	G	469	ILE	2.2
1	J	326	LYS	2.2
1	G	201	PHE	2.2
3	F	187	TRP	2.2
3	F	233	PHE	2.2
2	E	134	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	179	ASP	2.2
1	D	365	ILE	2.2
2	B	82	SER	2.2
1	D	209	ALA	2.2
1	G	319	ASN	2.2
1	J	287	ASN	2.2
1	J	411	PRO	2.2
3	F	131	LYS	2.2
3	I	92	GLY	2.2
1	D	221	ILE	2.1
1	A	266	SER	2.1
1	J	190	SER	2.1
2	H	68	ALA	2.1
1	J	198	ASN	2.1
2	E	33	PRO	2.1
1	G	452	LYS	2.1
2	H	39	LEU	2.1
3	F	127	LYS	2.1
2	E	138	ALA	2.1
1	A	181	SER	2.1
1	J	299	ILE	2.1
1	J	461	LEU	2.1
1	G	414	VAL	2.1
1	D	449	HIS	2.1
1	G	189	ILE	2.1
2	E	198	LEU	2.1
1	A	207	ARG	2.1
3	F	35	MET	2.1
1	D	463	GLY	2.1
1	A	424	PHE	2.1
1	G	462	PHE	2.1
3	F	34	PHE	2.1
1	D	445	LYS	2.1
1	G	372	ASN	2.1
1	A	405	LEU	2.1
1	G	285	LEU	2.1
3	F	71	LEU	2.1
3	I	97	LEU	2.1
1	G	280	ALA	2.1
2	E	111	ALA	2.1
1	G	317	GLN	2.1
1	G	364	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
2	E	146	TYR	2.0
1	A	462	PHE	2.0
3	I	86	PHE	2.0
1	A	451	GLU	2.0
1	J	219	GLU	2.0
1	J	410	LYS	2.0
1	J	307	LEU	2.0
2	E	203	SER	2.0
1	D	340	TYR	2.0
1	G	340	TYR	2.0
2	E	50	ASP	2.0
1	A	452	LYS	2.0
1	A	232	GLU	2.0
1	A	362	GLU	2.0
2	B	20	GLU	2.0
3	C	53	VAL	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.