



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:46 PM UTC

PDB ID : 4GTU / pdb_00004gtu
Title : LIGAND-FREE HOMODIMERIC HUMAN GLUTATHIONE S-TRANSFERASE M4-4
Authors : Patskovsky, Y.V.; Patskovska, L.N.; Listowsky, I.
Deposited on : 1999-05-12
Resolution : 3.30 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

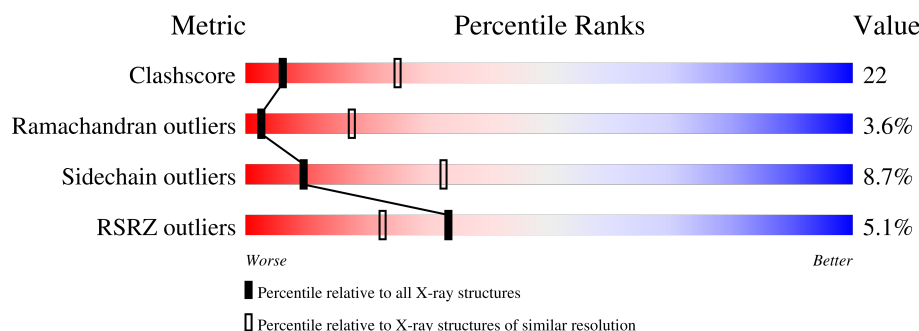
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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.



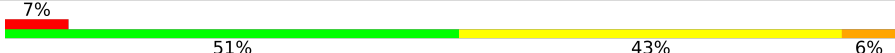
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	217	7% 50% 46% 5%
1	B	217	4% 53% 40% 8%
1	C	217	4% 47% 45% 7%
1	D	217	6% 50% 43% 7%
1	E	217	6% 53% 40% 6%
1	F	217	3% 39% 52% 9%

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Mol	Chain	Length	Quality of chain
1	G	217	 7% 51% 43% 6%
1	H	217	 5% 53% 41% 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 14352 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.

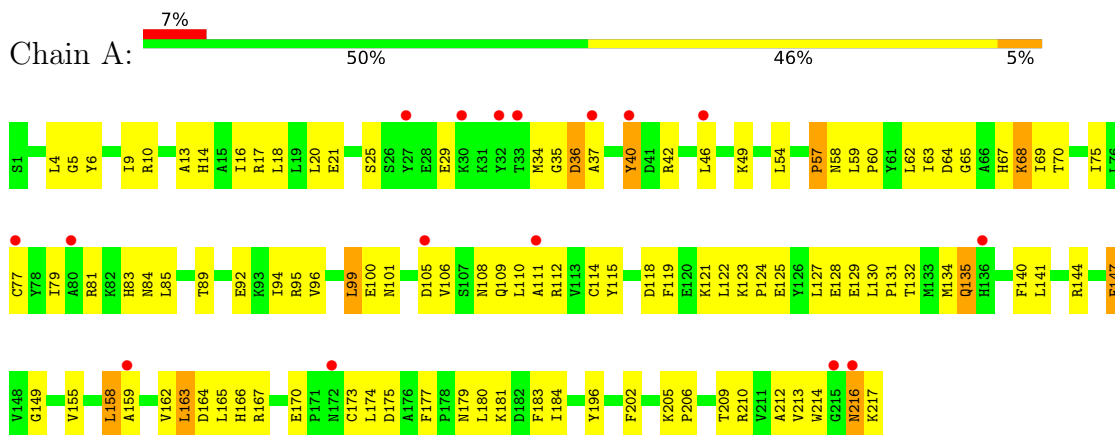
- Molecule 1 is a protein called GLUTATHIONE S-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	B	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	C	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	D	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	E	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	F	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	G	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			
1	H	217	Total	C	N	O	S	0	0	0
			1794	1160	295	329	10			

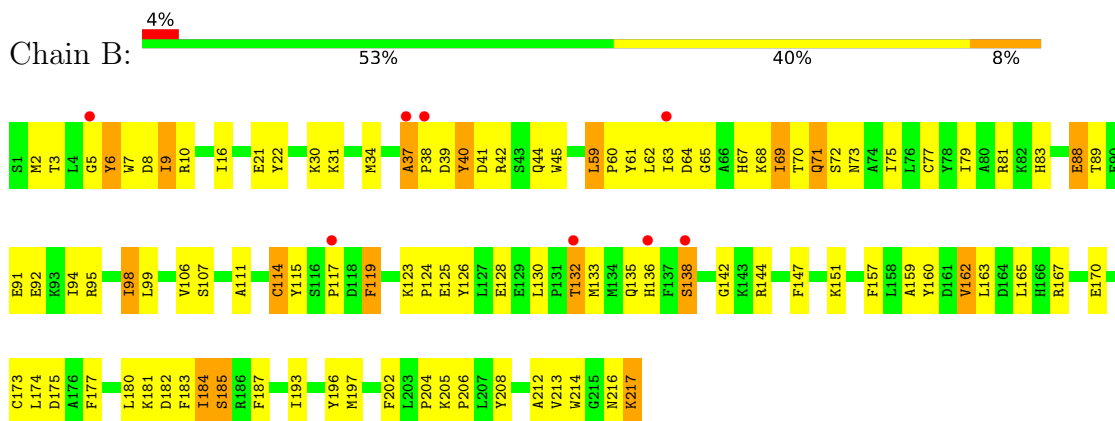
3 Residue-property plots

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.

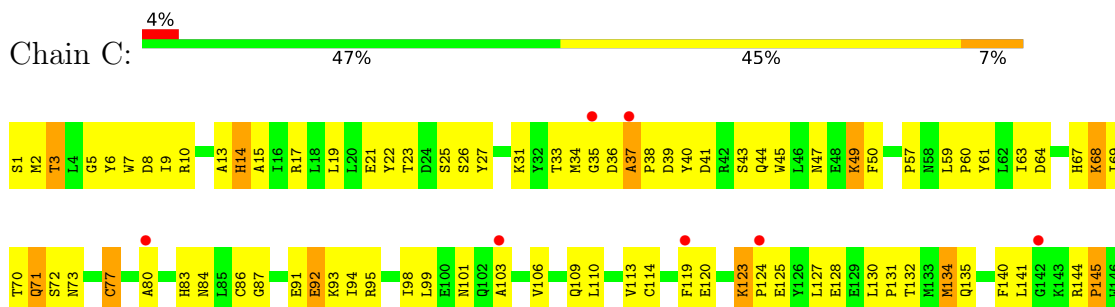
• Molecule 1: GLUTATHIONE S-TRANSFERASE



• Molecule 1: GLUTATHIONE S-TRANSFERASE

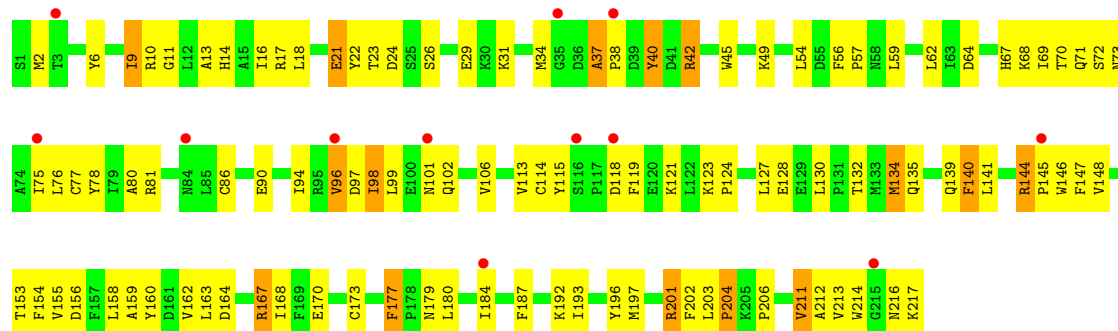


• Molecule 1: GLUTATHIONE S-TRANSFERASE

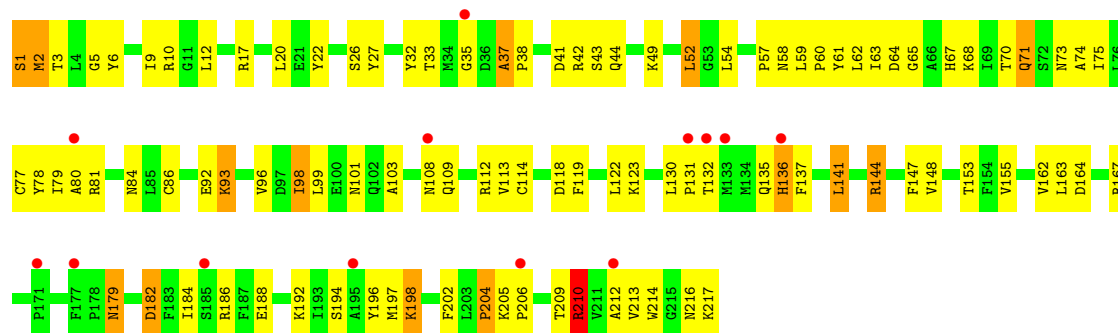




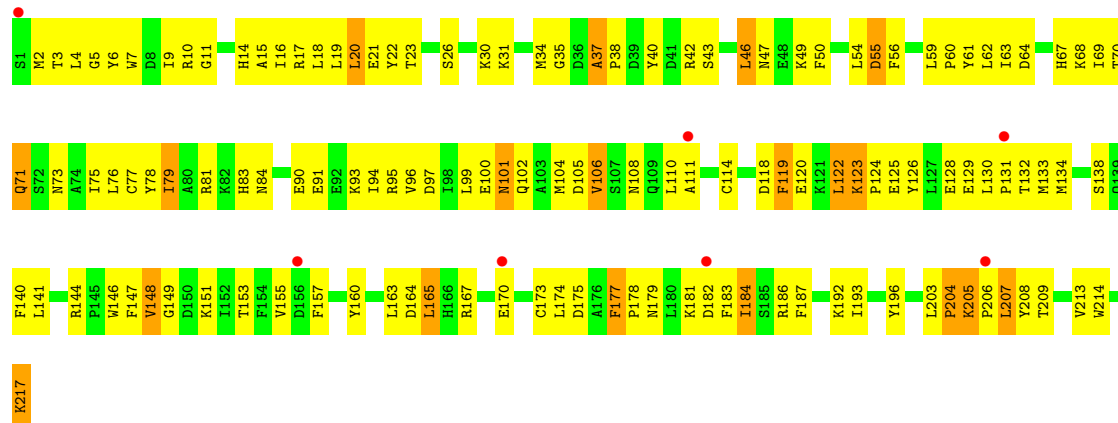
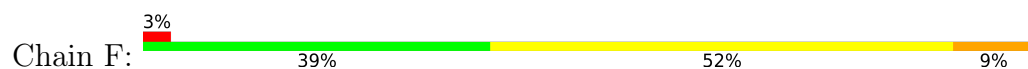
• Molecule 1: GLUTATHIONE S-TRANSFERASE



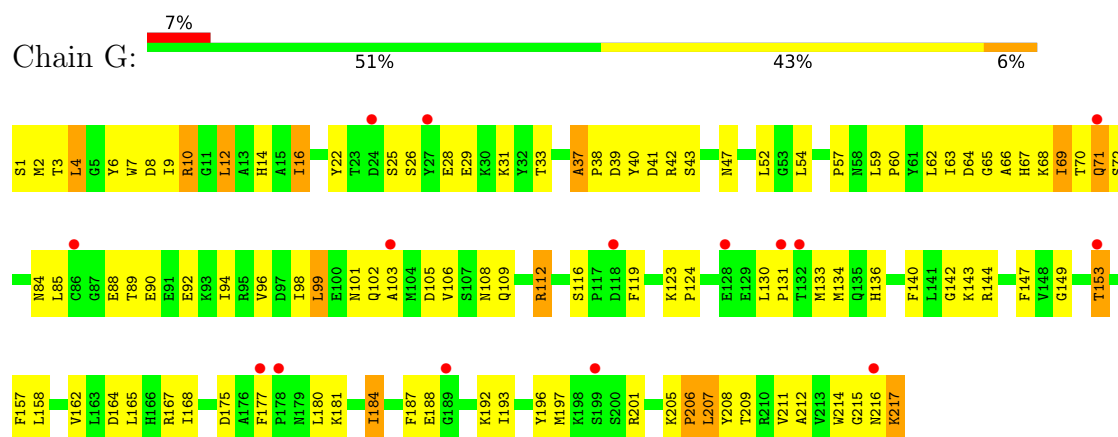
• Molecule 1: GLUTATHIONE S-TRANSFERASE



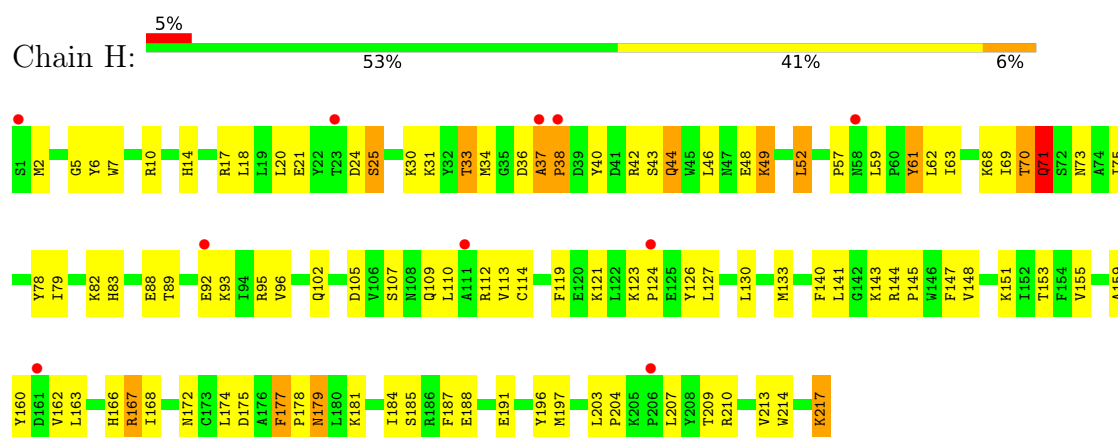
• Molecule 1: GLUTATHIONE S-TRANSFERASE



• Molecule 1: GLUTATHIONE S-TRANSFERASE



• Molecule 1: GLUTATHIONE S-TRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.00Å 215.70Å 94.56Å 90.00° 97.40° 90.00°	Depositor
Resolution (Å)	10.00 – 3.30 10.00 – 3.30	Depositor EDS
% Data completeness (in resolution range)	79.5 (10.00-3.30) 83.5 (10.00-3.30)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.81Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.245 , 0.315 0.259 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.16 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.77	EDS
Total number of atoms	14352	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.48	0/1842	0.93	3/2489 (0.1%)
1	B	0.51	0/1842	0.93	5/2489 (0.2%)
1	C	0.48	0/1842	0.94	3/2489 (0.1%)
1	D	0.48	0/1842	0.93	6/2489 (0.2%)
1	E	0.48	0/1842	0.96	1/2489 (0.0%)
1	F	0.48	0/1842	0.88	2/2489 (0.1%)
1	G	0.52	1/1842 (0.1%)	0.95	5/2489 (0.2%)
1	H	0.48	0/1842	0.93	3/2489 (0.1%)
All	All	0.49	1/14736 (0.0%)	0.93	28/19912 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	72	SER	C-N	-7.13	1.24	1.33

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	144	ARG	CA-C-N	7.37	126.88	118.85
1	D	144	ARG	C-N-CA	7.37	126.88	118.85
1	G	29	GLU	N-CA-C	6.01	118.52	108.96
1	B	177	PHE	CA-C-N	5.85	125.53	119.56
1	B	177	PHE	C-N-CA	5.85	125.53	119.56
1	C	177	PHE	CA-C-N	5.80	125.50	119.82
1	C	177	PHE	C-N-CA	5.80	125.50	119.82
1	A	110	LEU	N-CA-C	-5.77	105.13	111.82
1	E	141	LEU	N-CA-C	-5.59	104.82	111.03
1	B	88	GLU	N-CA-C	-5.54	104.69	112.12
1	B	107	SER	N-CA-C	-5.51	105.19	111.14
1	H	172	ASN	N-CA-C	5.51	119.26	112.54
1	C	179	ASN	N-CA-C	-5.39	104.44	111.02
1	H	177	PHE	CA-C-N	5.37	124.83	119.24
1	H	177	PHE	C-N-CA	5.37	124.83	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	GLU	N-CA-C	5.35	117.96	109.24
1	D	140	PHE	N-CA-C	-5.29	105.89	112.93
1	D	96	VAL	N-CA-C	-5.20	105.47	110.72
1	G	116	SER	CA-C-N	5.16	124.78	119.05
1	G	116	SER	C-N-CA	5.16	124.78	119.05
1	B	185	SER	N-CA-C	-5.16	106.07	112.93
1	A	112	ARG	N-CA-C	-5.12	105.62	111.14
1	D	177	PHE	CA-C-N	5.08	126.19	119.84
1	D	177	PHE	C-N-CA	5.08	126.19	119.84
1	F	177	PHE	CA-C-N	5.06	125.09	119.32
1	F	177	PHE	C-N-CA	5.06	125.09	119.32
1	G	168	ILE	N-CA-C	-5.06	105.22	112.35
1	G	69	ILE	O-C-N	5.00	128.13	122.67

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	1794	0	1762	85	0
1	B	1794	0	1762	85	0
1	C	1794	0	1762	87	0
1	D	1794	0	1762	77	0
1	E	1794	0	1762	76	0
1	F	1794	0	1762	109	0
1	G	1794	0	1762	75	0
1	H	1794	0	1762	73	0
All	All	14352	0	14096	639	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 (639) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:PRO:HG3	1:C:217:LYS:HB3	1.50	0.94
1:A:144:ARG:HD3	1:A:149:GLY:HA2	1.50	0.92
1:C:37:ALA:HB1	1:C:38:PRO:HD2	1.55	0.89
1:E:80:ALA:HB1	1:E:86:CYS:SG	2.12	0.89
1:F:134:MET:HE1	1:F:174:LEU:HA	1.56	0.88
1:A:114:CYS:SG	1:A:214:TRP:HB3	2.14	0.87
1:G:4:LEU:HD13	1:G:16:ILE:HG21	1.60	0.83
1:B:114:CYS:HG	1:B:214:TRP:HE3	1.24	0.83
1:A:63:ILE:HG12	1:A:68:LYS:HG3	1.62	0.82
1:G:1:SER:HB2	1:G:26:SER:O	1.79	0.81
1:H:174:LEU:HG	1:H:181:LYS:HD2	1.62	0.81
1:H:5:GLY:HA3	1:H:61:TYR:CE1	2.16	0.81
1:F:62:LEU:HD22	1:F:75:ILE:HG23	1.62	0.81
1:H:37:ALA:HB1	1:H:38:PRO:HD2	1.61	0.80
1:E:6:TYR:HA	1:E:59:LEU:HD22	1.63	0.80
1:E:71:GLN:HG2	1:F:101:ASN:HB2	1.62	0.80
1:E:210:ARG:HA	1:E:216:ASN:HB2	1.63	0.80
1:D:37:ALA:HB1	1:D:38:PRO:HD2	1.64	0.78
1:B:37:ALA:HB1	1:B:38:PRO:HD2	1.66	0.78
1:F:21:GLU:HG2	1:F:196:TYR:HB2	1.66	0.78
1:B:2:MET:HE2	1:B:62:LEU:HD11	1.66	0.78
1:G:101:ASN:HB3	1:H:71:GLN:HG2	1.66	0.78
1:F:9:ILE:HG13	1:F:206:PRO:HG2	1.65	0.77
1:F:34:MET:HE2	1:F:42:ARG:HG2	1.65	0.77
1:A:164:ASP:HA	1:A:167:ARG:HG3	1.68	0.76
1:H:6:TYR:HA	1:H:59:LEU:HD22	1.66	0.76
1:A:36:ASP:HA	1:A:40:TYR:CD2	2.21	0.75
1:D:115:TYR:HE1	1:D:212:ALA:HB2	1.52	0.75
1:H:174:LEU:HD21	1:H:184:ILE:HD12	1.66	0.75
1:E:37:ALA:HB1	1:E:38:PRO:HD2	1.68	0.74
1:E:136:HIS:HB3	1:F:55:ASP:O	1.85	0.74
1:G:37:ALA:HB1	1:G:38:PRO:HD2	1.69	0.74
1:B:2:MET:HB2	1:B:63:ILE:O	1.88	0.74
1:B:9:ILE:HG13	1:B:206:PRO:HG2	1.71	0.71
1:E:99:LEU:HD11	1:E:147:PHE:HB3	1.73	0.71
1:F:217:LYS:HB2	1:F:217:LYS:HZ2	1.55	0.71
1:F:90:GLU:HA	1:F:93:LYS:HD2	1.72	0.71
1:C:6:TYR:HA	1:C:59:LEU:HD22	1.71	0.70
1:A:135:GLN:HG2	1:A:177:PHE:CZ	2.25	0.70
1:B:175:ASP:HA	1:B:181:LYS:HD2	1.73	0.70
1:C:110:LEU:O	1:C:113:VAL:HG22	1.91	0.70
1:D:9:ILE:HG13	1:D:206:PRO:HG2	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:SER:HA	1:C:26:SER:O	1.92	0.70
1:C:144:ARG:HD3	1:C:149:GLY:HA2	1.72	0.70
1:D:158:LEU:O	1:D:162:VAL:HG23	1.92	0.70
1:D:160:TYR:HE1	1:D:184:ILE:HG23	1.55	0.70
1:F:6:TYR:CD1	1:F:60:PRO:HG3	2.27	0.70
1:G:164:ASP:HA	1:G:167:ARG:NH1	2.07	0.70
1:B:114:CYS:SG	1:B:214:TRP:HE3	2.15	0.69
1:D:206:PRO:HB3	1:D:217:LYS:HB2	1.73	0.69
1:G:205:LYS:HG3	1:G:217:LYS:HE2	1.74	0.69
1:D:216:ASN:O	1:D:217:LYS:HG3	1.93	0.69
1:C:134:MET:SD	1:C:174:LEU:HD12	2.33	0.69
1:D:6:TYR:O	1:D:31:LYS:HA	1.93	0.68
1:D:167:ARG:HG2	1:D:168:ILE:HD12	1.74	0.68
1:F:54:LEU:HD23	1:F:68:LYS:HB3	1.74	0.68
1:C:2:MET:HB2	1:C:63:ILE:O	1.93	0.68
1:C:110:LEU:HD22	1:C:166:HIS:CE1	2.29	0.68
1:C:130:LEU:HB3	1:C:131:PRO:HD3	1.75	0.68
1:F:163:LEU:HD22	1:F:174:LEU:HD11	1.76	0.68
1:C:99:LEU:HD11	1:C:147:PHE:HB3	1.77	0.68
1:D:123:LYS:HB3	1:D:124:PRO:HD3	1.76	0.68
1:A:64:ASP:HB3	1:A:67:HIS:HB2	1.76	0.67
1:F:110:LEU:HD23	1:F:165:LEU:HD12	1.77	0.67
1:D:29:GLU:HG2	1:D:31:LYS:HE3	1.76	0.67
1:D:96:VAL:HG13	1:D:155:VAL:HG13	1.77	0.67
1:F:73:ASN:HA	1:F:76:LEU:HD12	1.76	0.67
1:F:134:MET:SD	1:F:173:CYS:HB2	2.35	0.67
1:A:167:ARG:HD3	1:A:174:LEU:HD22	1.78	0.66
1:B:114:CYS:SG	1:B:214:TRP:HB3	2.35	0.66
1:C:164:ASP:O	1:C:167:ARG:HG2	1.94	0.66
1:A:114:CYS:HG	1:A:214:TRP:HE3	1.42	0.66
1:A:140:PHE:HD2	1:A:147:PHE:HD2	1.44	0.66
1:D:140:PHE:HD2	1:D:147:PHE:HD2	1.44	0.66
1:H:5:GLY:HA3	1:H:61:TYR:HE1	1.58	0.66
1:E:194:SER:O	1:E:198:LYS:HG2	1.94	0.66
1:E:62:LEU:HD22	1:E:75:ILE:HG23	1.78	0.65
1:C:95:ARG:NH2	1:C:144:ARG:HH12	1.95	0.65
1:F:9:ILE:HG12	1:F:209:THR:HA	1.78	0.65
1:A:101:ASN:HB3	1:B:71:GLN:OE1	1.96	0.65
1:E:78:TYR:HA	1:E:81:ARG:CZ	2.26	0.65
1:F:37:ALA:HB1	1:F:38:PRO:HD2	1.77	0.65
1:B:159:ALA:O	1:B:163:LEU:HG	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:VAL:HG11	1:D:153:THR:HG23	1.78	0.65
1:C:103:ALA:HB1	1:C:162:VAL:HG21	1.78	0.64
1:D:34:MET:HE2	1:D:42:ARG:HG2	1.78	0.64
1:D:160:TYR:CE1	1:D:184:ILE:HG23	2.32	0.64
1:G:205:LYS:CG	1:G:217:LYS:HE2	2.28	0.64
1:C:3:THR:HG23	1:C:63:ILE:HB	1.79	0.64
1:C:70:THR:O	1:C:71:GLN:HB2	1.98	0.64
1:H:88:GLU:HG2	1:H:92:GLU:OE2	1.98	0.64
1:H:123:LYS:HB3	1:H:124:PRO:HD3	1.79	0.64
1:B:77:CYS:HB3	1:B:81:ARG:HH12	1.61	0.63
1:E:167:ARG:CZ	1:E:184:ILE:HG21	2.28	0.63
1:F:129:GLU:HB2	1:F:133:MET:HE2	1.79	0.63
1:H:33:THR:HG22	1:H:210:ARG:HH21	1.63	0.63
1:B:167:ARG:NH2	1:B:184:ILE:HG21	2.12	0.63
1:C:119:PHE:HB2	1:C:213:VAL:HG21	1.79	0.63
1:C:164:ASP:HA	1:C:167:ARG:HD3	1.81	0.63
1:H:79:ILE:O	1:H:83:HIS:HD2	1.82	0.63
1:C:205:LYS:HD3	1:C:217:LYS:NZ	2.14	0.63
1:B:62:LEU:O	1:B:68:LYS:HA	1.99	0.63
1:E:114:CYS:HG	1:E:214:TRP:HE3	1.46	0.63
1:D:135:GLN:HG2	1:D:139:GLN:OE1	1.99	0.62
1:F:130:LEU:HB3	1:F:131:PRO:HD3	1.81	0.62
1:F:10:ARG:HD2	1:F:204:PRO:O	1.99	0.62
1:H:2:MET:SD	1:H:62:LEU:HD21	2.40	0.62
1:G:7:TRP:HH2	1:G:42:ARG:HG3	1.64	0.62
1:H:63:ILE:HD12	1:H:68:LYS:HG2	1.82	0.62
1:E:73:ASN:HB2	1:F:101:ASN:OD1	2.00	0.61
1:E:101:ASN:OD1	1:F:73:ASN:HB2	1.99	0.61
1:F:67:HIS:HB3	1:F:69:ILE:HD11	1.81	0.61
1:F:111:ALA:HA	1:F:208:TYR:HE2	1.64	0.61
1:E:141:LEU:HD13	1:E:147:PHE:CD1	2.36	0.61
1:D:170:GLU:HB3	1:D:173:CYS:SG	2.40	0.61
1:E:9:ILE:HG12	1:E:209:THR:HA	1.81	0.61
1:D:99:LEU:HD12	1:D:155:VAL:HB	1.81	0.61
1:F:2:MET:HA	1:F:63:ILE:O	2.00	0.61
1:B:21:GLU:HG2	1:B:196:TYR:HB2	1.83	0.61
1:E:1:SER:HA	1:E:26:SER:O	2.00	0.61
1:G:105:ASP:O	1:G:109:GLN:HG3	2.01	0.61
1:H:119:PHE:HB2	1:H:213:VAL:HG21	1.83	0.61
1:B:157:PHE:CD2	1:B:187:PHE:HZ	2.18	0.61
1:D:10:ARG:HG3	1:D:14:HIS:HB2	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:LEU:O	1:A:57:PRO:HD3	2.01	0.61
1:G:144:ARG:H	1:G:144:ARG:HD2	1.64	0.61
1:C:141:LEU:O	1:C:144:ARG:HG2	2.01	0.60
1:G:2:MET:CE	1:G:25:SER:HB3	2.31	0.60
1:B:205:LYS:HB3	1:B:206:PRO:HA	1.82	0.60
1:F:207:LEU:HD12	1:F:208:TYR:HE1	1.65	0.60
1:C:80:ALA:HB1	1:C:86:CYS:SG	2.41	0.60
1:E:184:ILE:O	1:E:188:GLU:HG2	2.01	0.60
1:F:6:TYR:O	1:F:31:LYS:HA	2.01	0.60
1:H:160:TYR:HE1	1:H:184:ILE:HG23	1.67	0.60
1:A:101:ASN:OD1	1:B:73:ASN:HB2	2.01	0.60
1:D:64:ASP:HB3	1:D:67:HIS:HB2	1.83	0.60
1:F:6:TYR:HA	1:F:59:LEU:HD22	1.83	0.59
1:F:207:LEU:HD12	1:F:208:TYR:CE1	2.37	0.59
1:E:153:THR:OG1	1:E:155:VAL:HG22	2.02	0.59
1:H:95:ARG:HB2	1:H:148:VAL:HG23	1.85	0.59
1:C:2:MET:HE1	1:C:25:SER:HB3	1.84	0.59
1:H:20:LEU:HD21	1:H:79:ILE:HD13	1.84	0.59
1:D:31:LYS:HE2	1:D:203:LEU:HD11	1.83	0.59
1:E:10:ARG:HB3	1:E:206:PRO:O	2.03	0.59
1:E:35:GLY:O	1:E:210:ARG:HG3	2.02	0.59
1:A:162:VAL:O	1:A:166:HIS:HD2	1.85	0.59
1:D:134:MET:HG2	1:D:173:CYS:HB3	1.83	0.59
1:H:112:ARG:HH11	1:H:112:ARG:HG2	1.68	0.59
1:A:4:LEU:HD12	1:A:20:LEU:HD21	1.85	0.59
1:D:6:TYR:HA	1:D:59:LEU:HD22	1.83	0.59
1:H:34:MET:HE2	1:H:42:ARG:HG2	1.84	0.59
1:G:144:ARG:HG2	1:G:149:GLY:HA2	1.84	0.59
1:C:10:ARG:HA	1:C:17:ARG:HH12	1.66	0.58
1:H:17:ARG:NH1	1:H:203:LEU:HD23	2.18	0.58
1:C:36:ASP:HA	1:C:40:TYR:HD1	1.67	0.58
1:D:115:TYR:CE1	1:D:212:ALA:HB2	2.34	0.58
1:G:9:ILE:HG13	1:G:209:THR:HA	1.85	0.58
1:B:160:TYR:CE1	1:B:184:ILE:HG23	2.38	0.58
1:C:9:ILE:HG12	1:C:209:THR:HA	1.86	0.58
1:E:20:LEU:HD23	1:E:79:ILE:HD11	1.84	0.58
1:G:103:ALA:O	1:G:106:VAL:HG22	2.03	0.58
1:B:40:TYR:O	1:B:42:ARG:HG2	2.04	0.58
1:H:21:GLU:HG2	1:H:196:TYR:HB2	1.85	0.58
1:B:119:PHE:HB2	1:B:213:VAL:HG11	1.84	0.58
1:D:119:PHE:HB2	1:D:213:VAL:HG11	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:217:LYS:HB2	1:F:217:LYS:NZ	2.17	0.58
1:A:9:ILE:HG13	1:A:209:THR:HA	1.85	0.58
1:G:63:ILE:HA	1:G:67:HIS:O	2.04	0.58
1:B:114:CYS:O	1:B:213:VAL:HG12	2.04	0.57
1:E:78:TYR:HA	1:E:81:ARG:NH1	2.19	0.57
1:A:114:CYS:O	1:A:213:VAL:HG22	2.03	0.57
1:E:206:PRO:HG3	1:E:217:LYS:HB2	1.85	0.57
1:H:18:LEU:HD21	1:H:187:PHE:HE2	1.69	0.57
1:A:163:LEU:HD23	1:A:184:ILE:HD11	1.87	0.57
1:E:3:THR:HG22	1:E:63:ILE:HB	1.86	0.57
1:H:37:ALA:HB1	1:H:38:PRO:CD	2.34	0.57
1:A:6:TYR:HA	1:A:59:LEU:HG	1.87	0.57
1:F:140:PHE:O	1:F:144:ARG:NH1	2.38	0.57
1:A:205:LYS:HG3	1:A:206:PRO:HA	1.87	0.57
1:B:206:PRO:HB3	1:B:217:LYS:CB	2.34	0.57
1:F:3:THR:O	1:F:62:LEU:HA	2.04	0.57
1:A:6:TYR:HD2	1:A:17:ARG:HH22	1.53	0.56
1:A:175:ASP:HA	1:A:181:LYS:CE	2.35	0.56
1:C:5:GLY:HA3	1:C:61:TYR:CE2	2.40	0.56
1:G:63:ILE:HG12	1:G:68:LYS:HG2	1.87	0.56
1:C:35:GLY:HA3	1:C:41:ASP:HB2	1.87	0.56
1:A:175:ASP:HA	1:A:181:LYS:HE2	1.87	0.56
1:D:164:ASP:HA	1:D:167:ARG:HD2	1.87	0.56
1:F:22:TYR:HA	1:F:192:LYS:HB3	1.88	0.56
1:C:6:TYR:CD1	1:C:60:PRO:HG3	2.40	0.56
1:F:21:GLU:HG2	1:F:196:TYR:CB	2.35	0.56
1:F:175:ASP:HA	1:F:181:LYS:HD3	1.88	0.56
1:C:110:LEU:HD22	1:C:166:HIS:HE1	1.70	0.56
1:C:210:ARG:HA	1:C:216:ASN:HB2	1.87	0.56
1:G:64:ASP:HB3	1:G:67:HIS:HB2	1.88	0.55
1:E:119:PHE:HB2	1:E:213:VAL:HG21	1.87	0.55
1:C:43:SER:O	1:C:47:ASN:HB2	2.06	0.55
1:D:13:ALA:O	1:D:17:ARG:HG3	2.05	0.55
1:C:23:THR:HB	1:C:83:HIS:CD2	2.42	0.55
1:D:114:CYS:HG	1:D:214:TRP:HE3	1.51	0.55
1:F:110:LEU:HD13	1:F:126:TYR:CE2	2.41	0.55
1:B:45:TRP:CH2	1:B:61:TYR:CE2	2.94	0.55
1:C:71:GLN:HG3	1:D:101:ASN:HB3	1.88	0.55
1:E:17:ARG:HA	1:E:20:LEU:HD12	1.89	0.55
1:D:14:HIS:O	1:D:18:LEU:HG	2.06	0.55
1:A:10:ARG:HG3	1:A:14:HIS:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:136:HIS:CD2	1:H:57:PRO:HD2	2.42	0.55
1:G:196:TYR:HE1	1:G:201:ARG:HB2	1.72	0.55
1:F:15:ALA:HA	1:F:18:LEU:HD12	1.89	0.55
1:A:54:LEU:HD23	1:A:68:LYS:HB3	1.88	0.54
1:A:130:LEU:O	1:A:134:MET:HG2	2.07	0.54
1:E:70:THR:O	1:E:71:GLN:HB2	2.07	0.54
1:F:77:CYS:O	1:F:81:ARG:HG3	2.07	0.54
1:A:114:CYS:HG	1:A:214:TRP:HB3	1.73	0.54
1:F:114:CYS:O	1:F:213:VAL:HG22	2.07	0.54
1:G:4:LEU:CD1	1:G:16:ILE:HG21	2.36	0.54
1:E:92:GLU:O	1:E:96:VAL:HG23	2.08	0.54
1:G:207:LEU:HD12	1:G:208:TYR:CE1	2.42	0.54
1:H:93:LYS:O	1:H:96:VAL:HB	2.07	0.54
1:B:184:ILE:HG22	1:B:184:ILE:O	2.08	0.54
1:G:70:THR:O	1:G:71:GLN:HB2	2.07	0.54
1:D:90:GLU:HG3	1:D:94:ILE:HD11	1.90	0.54
1:E:20:LEU:HD23	1:E:79:ILE:CD1	2.37	0.54
1:A:54:LEU:HD12	1:A:57:PRO:HA	1.89	0.53
1:D:130:LEU:HD21	1:D:173:CYS:SG	2.47	0.53
1:E:93:LYS:HG2	1:F:81:ARG:HE	1.72	0.53
1:F:78:TYR:O	1:F:81:ARG:HB2	2.08	0.53
1:G:12:LEU:HD13	1:G:60:PRO:HD3	1.88	0.53
1:C:10:ARG:HG3	1:C:14:HIS:HB2	1.90	0.53
1:F:129:GLU:O	1:F:133:MET:HG3	2.09	0.53
1:A:16:ILE:O	1:A:20:LEU:HG	2.09	0.53
1:A:83:HIS:O	1:A:85:LEU:HD12	2.08	0.53
1:C:114:CYS:O	1:C:213:VAL:HG22	2.09	0.53
1:A:173:CYS:SG	1:A:174:LEU:N	2.81	0.53
1:D:114:CYS:SG	1:D:214:TRP:CE3	3.01	0.53
1:E:136:HIS:HB2	1:F:56:PHE:CE1	2.44	0.53
1:A:63:ILE:HA	1:A:67:HIS:O	2.09	0.53
1:E:77:CYS:O	1:E:81:ARG:HG2	2.09	0.53
1:E:148:VAL:HG21	1:E:153:THR:OG1	2.09	0.53
1:A:135:GLN:HG2	1:A:177:PHE:HZ	1.74	0.53
1:F:10:ARG:HE	1:F:14:HIS:CG	2.27	0.53
1:F:71:GLN:OE1	1:F:71:GLN:HA	2.08	0.53
1:A:167:ARG:NH1	1:A:184:ILE:HD13	2.24	0.53
1:B:77:CYS:HB3	1:B:81:ARG:NH1	2.23	0.53
1:B:183:PHE:C	1:B:185:SER:H	2.17	0.53
1:H:48:GLU:O	1:H:52:LEU:HD12	2.09	0.52
1:F:21:GLU:HG2	1:F:196:TYR:CG	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ARG:HG3	1:A:46:LEU:HD11	1.91	0.52
1:F:134:MET:SD	1:F:174:LEU:HD22	2.50	0.52
1:H:110:LEU:O	1:H:113:VAL:HG22	2.09	0.52
1:B:217:LYS:HB2	1:B:217:LYS:NZ	2.23	0.52
1:G:177:PHE:HB2	1:G:180:LEU:HD12	1.91	0.52
1:C:45:TRP:CZ2	1:C:49:LYS:HG3	2.44	0.52
1:F:160:TYR:CD1	1:F:184:ILE:HG22	2.45	0.52
1:G:69:ILE:HD12	1:G:69:ILE:H	1.74	0.52
1:D:202:PHE:O	1:D:204:PRO:HD3	2.10	0.52
1:A:196:TYR:CE2	1:A:202:PHE:HD1	2.27	0.52
1:B:91:GLU:O	1:B:95:ARG:HG3	2.09	0.52
1:G:9:ILE:CD1	1:G:206:PRO:HB3	2.40	0.52
1:G:101:ASN:OD1	1:H:73:ASN:HB2	2.10	0.52
1:G:205:LYS:HA	1:G:206:PRO:C	2.35	0.52
1:A:123:LYS:HB3	1:A:124:PRO:HD3	1.91	0.52
1:B:5:GLY:HA2	1:B:30:LYS:O	2.10	0.52
1:F:4:LEU:HD21	1:F:17:ARG:HG2	1.91	0.52
1:F:102:GLN:O	1:F:106:VAL:HG23	2.10	0.52
1:H:217:LYS:HB2	1:H:217:LYS:NZ	2.24	0.52
1:F:30:LYS:HE3	1:F:61:TYR:OH	2.10	0.52
1:B:64:ASP:HB3	1:B:67:HIS:HB2	1.90	0.51
1:D:96:VAL:HG13	1:D:155:VAL:CG1	2.39	0.51
1:E:54:LEU:HD23	1:E:68:LYS:HB3	1.92	0.51
1:F:111:ALA:HA	1:F:208:TYR:CE2	2.44	0.51
1:C:10:ARG:HD2	1:C:204:PRO:O	2.10	0.51
1:D:14:HIS:HE1	1:D:202:PHE:CE1	2.28	0.51
1:B:16:ILE:HD12	1:B:16:ILE:H	1.75	0.51
1:D:73:ASN:HA	1:D:76:LEU:HD12	1.93	0.51
1:A:94:ILE:HG23	1:B:69:ILE:HG21	1.93	0.51
1:D:21:GLU:HG3	1:D:196:TYR:CD1	2.45	0.51
1:F:35:GLY:O	1:F:40:TYR:HA	2.10	0.51
1:E:164:ASP:O	1:E:167:ARG:HB3	2.10	0.51
1:G:164:ASP:HA	1:G:167:ARG:HH11	1.73	0.51
1:A:118:ASP:HB3	1:A:121:LYS:HB2	1.92	0.51
1:H:179:ASN:N	1:H:179:ASN:HD22	2.09	0.51
1:C:164:ASP:HA	1:C:167:ARG:CD	2.40	0.51
1:E:6:TYR:HA	1:E:59:LEU:CD2	2.37	0.51
1:D:114:CYS:SG	1:D:214:TRP:HE3	2.33	0.51
1:E:6:TYR:CE1	1:E:60:PRO:HD3	2.46	0.51
1:H:126:TYR:CE1	1:H:133:MET:HE1	2.46	0.51
1:B:61:TYR:CE1	1:B:63:ILE:HD11	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:5:GLY:O	1:E:60:PRO:HA	2.11	0.50
1:B:88:GLU:CD	1:B:151:LYS:HE3	2.37	0.50
1:B:6:TYR:HA	1:B:59:LEU:HD22	1.93	0.50
1:C:36:ASP:O	1:C:37:ALA:HB3	2.11	0.50
1:F:62:LEU:CD2	1:F:69:ILE:HB	2.42	0.50
1:G:8:ASP:OD2	1:G:33:THR:HA	2.11	0.50
1:A:115:TYR:HE1	1:A:212:ALA:HB2	1.77	0.50
1:C:94:ILE:HG23	1:D:69:ILE:HD12	1.93	0.50
1:F:110:LEU:CD2	1:F:165:LEU:HD12	2.41	0.50
1:F:146:TRP:CH2	1:F:186:ARG:HG2	2.45	0.50
1:F:207:LEU:HD11	1:F:214:TRP:CZ3	2.46	0.50
1:G:85:LEU:O	1:G:153:THR:HG22	2.12	0.50
1:G:123:LYS:N	1:G:124:PRO:HD2	2.27	0.50
1:A:100:GLU:HA	1:A:158:LEU:HD13	1.94	0.50
1:D:77:CYS:O	1:D:81:ARG:HB2	2.10	0.50
1:E:210:ARG:HH11	1:E:210:ARG:HB3	1.77	0.50
1:G:130:LEU:O	1:G:133:MET:HB3	2.12	0.50
1:H:10:ARG:HG2	1:H:14:HIS:HB2	1.93	0.50
1:A:114:CYS:SG	1:A:214:TRP:HE3	2.35	0.50
1:B:216:ASN:O	1:B:217:LYS:HD3	2.12	0.50
1:F:22:TYR:HB2	1:F:193:ILE:HD11	1.93	0.50
1:D:78:TYR:HA	1:D:81:ARG:NH1	2.26	0.50
1:B:91:GLU:HG3	1:B:92:GLU:N	2.27	0.50
1:B:130:LEU:O	1:B:133:MET:HB3	2.12	0.50
1:C:69:ILE:HG23	1:D:98:ILE:HD11	1.94	0.50
1:G:119:PHE:CZ	1:G:214:TRP:CD1	3.00	0.50
1:B:6:TYR:CE1	1:B:60:PRO:HD3	2.47	0.49
1:B:193:ILE:HD12	1:B:193:ILE:N	2.27	0.49
1:E:58:ASN:HB3	1:E:70:THR:HG22	1.92	0.49
1:B:197:MET:SD	1:B:202:PHE:CZ	3.04	0.49
1:B:111:ALA:HA	1:B:208:TYR:CE2	2.47	0.49
1:D:187:PHE:CE2	1:D:193:ILE:HD13	2.46	0.49
1:F:23:THR:HB	1:F:83:HIS:CD2	2.47	0.49
1:B:111:ALA:HA	1:B:208:TYR:HE2	1.76	0.49
1:D:22:TYR:HD2	1:D:23:THR:HG23	1.76	0.49
1:E:10:ARG:HD2	1:E:204:PRO:O	2.12	0.49
1:F:205:LYS:HD3	1:F:217:LYS:CE	2.42	0.49
1:A:108:ASN:HA	1:A:111:ALA:HB3	1.94	0.49
1:C:73:ASN:HB2	1:D:101:ASN:OD1	2.12	0.49
1:H:187:PHE:CD2	1:H:187:PHE:O	2.65	0.49
1:A:18:LEU:O	1:A:21:GLU:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:CYS:O	1:C:80:ALA:HB3	2.13	0.49
1:B:59:LEU:HA	1:B:60:PRO:C	2.38	0.49
1:B:170:GLU:O	1:B:173:CYS:SG	2.66	0.49
1:F:205:LYS:HB2	1:F:206:PRO:HA	1.95	0.49
1:H:5:GLY:HA2	1:H:30:LYS:O	2.13	0.49
1:D:14:HIS:HE1	1:D:202:PHE:HE1	1.60	0.49
1:G:3:THR:HB	1:G:28:GLU:HB3	1.94	0.49
1:C:7:TRP:C	1:C:9:ILE:H	2.21	0.48
1:H:20:LEU:CD2	1:H:79:ILE:HD13	2.42	0.48
1:F:91:GLU:O	1:F:95:ARG:HG3	2.13	0.48
1:B:89:THR:OG1	1:B:92:GLU:HG3	2.13	0.48
1:B:174:LEU:H	1:B:174:LEU:HD12	1.78	0.48
1:D:124:PRO:O	1:D:128:GLU:HG3	2.14	0.48
1:D:78:TYR:HA	1:D:81:ARG:HH12	1.78	0.48
1:G:10:ARG:HG2	1:G:14:HIS:HB2	1.94	0.48
1:C:36:ASP:O	1:C:40:TYR:HA	2.14	0.48
1:F:205:LYS:HD3	1:F:217:LYS:HE2	1.94	0.48
1:G:88:GLU:HB2	1:G:92:GLU:OE1	2.13	0.48
1:G:212:ALA:O	1:G:216:ASN:ND2	2.46	0.48
1:E:119:PHE:CZ	1:E:123:LYS:HD3	2.48	0.48
1:G:69:ILE:HD12	1:G:69:ILE:N	2.28	0.48
1:D:102:GLN:O	1:D:106:VAL:HG23	2.14	0.48
1:D:148:VAL:HG11	1:D:153:THR:CG2	2.42	0.48
1:F:62:LEU:HD23	1:F:69:ILE:HB	1.96	0.48
1:F:119:PHE:HB2	1:F:213:VAL:HG21	1.95	0.48
1:C:25:SER:O	1:C:27:TYR:HD1	1.97	0.48
1:C:63:ILE:HG23	1:C:68:LYS:HE2	1.96	0.48
1:C:93:LYS:HD2	1:D:81:ARG:HD3	1.94	0.48
1:G:103:ALA:HB1	1:G:162:VAL:HG21	1.96	0.48
1:A:114:CYS:SG	1:A:214:TRP:CE3	3.07	0.47
1:C:123:LYS:HD3	1:C:127:LEU:HD11	1.95	0.47
1:D:118:ASP:HB3	1:D:121:LYS:HD2	1.95	0.47
1:G:9:ILE:HA	1:G:206:PRO:HB2	1.96	0.47
1:A:174:LEU:O	1:A:181:LYS:HE2	2.14	0.47
1:B:21:GLU:HG2	1:B:196:TYR:CG	2.49	0.47
1:A:180:LEU:O	1:A:183:PHE:HB3	2.15	0.47
1:B:8:ASP:HA	1:B:31:LYS:HD2	1.95	0.47
1:B:163:LEU:HB3	1:B:184:ILE:HD11	1.97	0.47
1:B:212:ALA:O	1:B:216:ASN:ND2	2.48	0.47
1:E:96:VAL:O	1:E:155:VAL:HG11	2.14	0.47
1:G:22:TYR:O	1:G:192:LYS:HD3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:MET:HG2	1:A:35:GLY:N	2.30	0.47
1:F:170:GLU:O	1:F:173:CYS:SG	2.64	0.47
1:B:123:LYS:N	1:B:124:PRO:HD2	2.30	0.47
1:A:141:LEU:HB3	1:A:179:ASN:HD21	1.80	0.47
1:A:167:ARG:HG3	1:A:167:ARG:HH11	1.80	0.47
1:B:132:THR:O	1:B:135:GLN:HB3	2.15	0.47
1:C:3:THR:HG23	1:C:63:ILE:HD12	1.97	0.47
1:C:49:LYS:HE2	1:C:57:PRO:HB2	1.96	0.47
1:C:181:LYS:O	1:C:184:ILE:HG22	2.14	0.47
1:A:118:ASP:O	1:A:122:LEU:HD13	2.15	0.47
1:D:159:ALA:O	1:D:163:LEU:HG	2.15	0.47
1:H:140:PHE:HD2	1:H:147:PHE:HD1	1.61	0.47
1:B:132:THR:O	1:B:136:HIS:ND1	2.47	0.47
1:E:70:THR:O	1:E:71:GLN:CB	2.62	0.47
1:A:105:ASP:O	1:A:109:GLN:HG3	2.15	0.47
1:A:115:TYR:CE1	1:A:212:ALA:HB2	2.50	0.46
1:B:6:TYR:O	1:B:31:LYS:HA	2.16	0.46
1:E:113:VAL:HG13	1:E:122:LEU:HB3	1.98	0.46
1:G:9:ILE:HD12	1:G:206:PRO:HB3	1.97	0.46
1:A:144:ARG:HD3	1:A:149:GLY:CA	2.34	0.46
1:F:43:SER:HA	1:F:46:LEU:HB2	1.97	0.46
1:E:22:TYR:HD1	1:E:192:LYS:HB2	1.80	0.46
1:F:157:PHE:CD2	1:F:187:PHE:HZ	2.33	0.46
1:C:13:ALA:O	1:C:15:ALA:N	2.49	0.46
1:D:40:TYR:CG	1:D:211:VAL:HG12	2.49	0.46
1:G:102:GLN:O	1:G:105:ASP:HB3	2.16	0.46
1:F:15:ALA:O	1:F:18:LEU:HB2	2.16	0.46
1:F:69:ILE:HD12	1:F:69:ILE:N	2.31	0.46
1:H:174:LEU:HG	1:H:181:LYS:CD	2.41	0.46
1:B:62:LEU:HD22	1:B:75:ILE:HG23	1.98	0.46
1:F:64:ASP:HB3	1:F:67:HIS:HB2	1.98	0.46
1:G:62:LEU:O	1:G:68:LYS:HA	2.16	0.46
1:B:9:ILE:HG13	1:B:206:PRO:CG	2.43	0.46
1:D:134:MET:HB3	1:D:177:PHE:CE2	2.51	0.46
1:H:33:THR:HG22	1:H:210:ARG:NH2	2.29	0.46
1:C:106:VAL:O	1:C:109:GLN:HB3	2.16	0.46
1:F:6:TYR:HE2	1:F:9:ILE:O	1.99	0.46
1:G:10:ARG:CG	1:G:14:HIS:HB2	2.46	0.46
1:G:99:LEU:HA	1:G:99:LEU:HD22	1.51	0.46
1:G:130:LEU:N	1:G:131:PRO:HD2	2.31	0.46
1:H:159:ALA:O	1:H:163:LEU:HG	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:207:LEU:N	1:H:207:LEU:HD22	2.31	0.46
1:E:130:LEU:N	1:E:131:PRO:HD2	2.31	0.45
1:E:196:TYR:O	1:E:202:PHE:HB2	2.16	0.45
1:F:110:LEU:HD13	1:F:126:TYR:HE2	1.81	0.45
1:F:134:MET:HE2	1:F:134:MET:HB3	1.69	0.45
1:F:141:LEU:HD21	1:F:183:PHE:HB2	1.97	0.45
1:G:94:ILE:HD13	1:H:69:ILE:HD11	1.97	0.45
1:A:205:LYS:HA	1:A:206:PRO:C	2.41	0.45
1:B:173:CYS:SG	1:B:174:LEU:HD12	2.57	0.45
1:H:107:SER:OG	1:H:162:VAL:HG22	2.17	0.45
1:A:129:GLU:O	1:A:132:THR:HG22	2.16	0.45
1:G:207:LEU:HD12	1:G:208:TYR:HE1	1.79	0.45
1:A:127:LEU:HD22	1:A:170:GLU:OE1	2.17	0.45
1:B:182:ASP:O	1:B:185:SER:HB2	2.17	0.45
1:C:209:THR:OG1	1:C:211:VAL:HG12	2.17	0.45
1:F:138:SER:O	1:F:179:ASN:ND2	2.49	0.45
1:H:43:SER:HA	1:H:46:LEU:HB2	1.99	0.45
1:A:106:VAL:HG13	1:A:162:VAL:HG11	1.99	0.45
1:B:7:TRP:C	1:B:9:ILE:H	2.24	0.45
1:B:22:TYR:CD1	1:B:193:ILE:HD11	2.51	0.45
1:B:65:GLY:C	1:B:67:HIS:H	2.25	0.45
1:E:179:ASN:HD22	1:E:179:ASN:N	2.15	0.45
1:A:99:LEU:HD12	1:A:155:VAL:HB	1.98	0.45
1:G:157:PHE:HA	1:G:187:PHE:CE2	2.52	0.45
1:E:80:ALA:CB	1:E:86:CYS:SG	2.95	0.45
1:C:95:ARG:NH1	1:C:140:PHE:CE2	2.85	0.45
1:F:148:VAL:HG12	1:F:151:LYS:O	2.15	0.45
1:G:71:GLN:HE21	1:H:102:GLN:HA	1.82	0.45
1:A:79:ILE:O	1:A:83:HIS:HD2	1.99	0.45
1:A:119:PHE:HB2	1:A:213:VAL:HG21	1.98	0.45
1:B:157:PHE:CE2	1:B:187:PHE:HZ	2.35	0.45
1:D:49:LYS:O	1:D:57:PRO:HB3	2.16	0.45
1:D:70:THR:O	1:D:71:GLN:HB2	2.16	0.45
1:B:5:GLY:HA2	1:B:30:LYS:HB3	1.98	0.44
1:C:49:LYS:HD3	1:C:50:PHE:CZ	2.52	0.44
1:F:22:TYR:O	1:F:192:LYS:HE3	2.16	0.44
1:A:77:CYS:O	1:A:81:ARG:HB2	2.17	0.44
1:B:125:GLU:O	1:B:128:GLU:HG2	2.17	0.44
1:E:167:ARG:NH1	1:E:184:ILE:HG21	2.32	0.44
1:C:5:GLY:HA3	1:C:61:TYR:CZ	2.52	0.44
1:D:54:LEU:HD12	1:D:57:PRO:HA	1.97	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:118:ASP:O	1:F:120:GLU:N	2.51	0.44
1:F:122:LEU:H	1:F:122:LEU:HD12	1.82	0.44
1:G:134:MET:HB2	1:G:177:PHE:CE2	2.51	0.44
1:B:106:VAL:HG23	1:B:162:VAL:HG21	1.98	0.44
1:G:2:MET:O	1:G:28:GLU:N	2.50	0.44
1:C:125:GLU:O	1:C:128:GLU:HG2	2.17	0.44
1:C:132:THR:O	1:C:135:GLN:HB3	2.18	0.44
1:F:144:ARG:HD3	1:F:149:GLY:HA2	2.00	0.44
1:C:101:ASN:HB3	1:D:71:GLN:OE1	2.18	0.44
1:G:140:PHE:HD2	1:G:147:PHE:HD1	1.66	0.44
1:G:193:ILE:HG22	1:G:197:MET:SD	2.57	0.44
1:H:7:TRP:HH2	1:H:42:ARG:HB3	1.82	0.44
1:C:8:ASP:HB3	1:C:31:LYS:HG2	1.98	0.44
1:C:70:THR:H	1:D:98:ILE:HD11	1.83	0.44
1:C:183:PHE:HA	1:C:186:ARG:NH1	2.33	0.44
1:A:62:LEU:HD22	1:A:75:ILE:HG23	2.00	0.44
1:E:78:TYR:HE1	1:F:94:ILE:HD11	1.83	0.44
1:H:177:PHE:HA	1:H:178:PRO:HD2	1.83	0.44
1:A:36:ASP:HA	1:A:40:TYR:CE2	2.52	0.44
1:B:117:PRO:C	1:B:119:PHE:H	2.26	0.44
1:D:86:CYS:O	1:D:153:THR:HG22	2.18	0.44
1:E:109:GLN:O	1:E:113:VAL:HG23	2.18	0.44
1:F:16:ILE:O	1:F:20:LEU:HB2	2.18	0.44
1:F:123:LYS:N	1:F:124:PRO:HD2	2.33	0.44
1:G:12:LEU:CD1	1:G:60:PRO:HD3	2.47	0.44
1:A:81:ARG:HH22	1:B:94:ILE:HG22	1.82	0.43
1:C:92:GLU:HG3	1:C:148:VAL:CG1	2.48	0.43
1:C:192:LYS:NZ	1:C:192:LYS:HB2	2.33	0.43
1:E:2:MET:O	1:E:27:TYR:HA	2.17	0.43
1:E:74:ALA:HA	1:F:97:ASP:HB3	1.99	0.43
1:E:163:LEU:HB2	1:E:184:ILE:HD11	2.00	0.43
1:F:5:GLY:HA3	1:F:61:TYR:CZ	2.53	0.43
1:G:184:ILE:HG23	1:G:188:GLU:OE2	2.17	0.43
1:C:196:TYR:HD1	1:C:201:ARG:HH21	1.65	0.43
1:F:164:ASP:O	1:F:167:ARG:HB3	2.18	0.43
1:H:24:ASP:O	1:H:25:SER:C	2.61	0.43
1:H:49:LYS:HD3	1:H:57:PRO:CB	2.48	0.43
1:H:160:TYR:CE1	1:H:184:ILE:HG23	2.51	0.43
1:H:167:ARG:HD3	1:H:168:ILE:HG13	2.01	0.43
1:A:216:ASN:HD22	1:A:216:ASN:H	1.66	0.43
1:H:153:THR:OG1	1:H:155:VAL:HG22	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:197:MET:HE3	1:H:197:MET:HB2	1.86	0.43
1:C:87:GLY:HA2	1:C:92:GLU:HB3	2.00	0.43
1:F:61:TYR:HB3	1:F:70:THR:HG22	2.00	0.43
1:G:207:LEU:HB3	1:G:208:TYR:HD1	1.82	0.43
1:A:58:ASN:CG	1:A:59:LEU:N	2.75	0.43
1:B:115:TYR:CE1	1:B:212:ALA:HB2	2.53	0.43
1:E:114:CYS:O	1:E:213:VAL:HG22	2.17	0.43
1:F:153:THR:OG1	1:F:155:VAL:HG22	2.18	0.43
1:H:143:LYS:HD3	1:H:143:LYS:N	2.34	0.43
1:C:21:GLU:HG3	1:C:196:TYR:CD1	2.54	0.43
1:C:71:GLN:CG	1:D:101:ASN:HB3	2.48	0.43
1:C:95:ARG:NH1	1:C:147:PHE:O	2.51	0.43
1:E:64:ASP:N	1:E:67:HIS:O	2.50	0.43
1:F:49:LYS:HD3	1:F:50:PHE:CE1	2.54	0.43
1:A:13:ALA:HA	1:A:16:ILE:HD12	2.00	0.43
1:C:205:LYS:HA	1:C:206:PRO:C	2.44	0.43
1:D:10:ARG:HG2	1:D:11:GLY:N	2.34	0.43
1:F:164:ASP:HA	1:F:167:ARG:CB	2.49	0.43
1:C:64:ASP:HB3	1:C:67:HIS:HB2	2.00	0.43
1:C:196:TYR:CZ	1:C:202:PHE:HA	2.54	0.43
1:E:10:ARG:HB2	1:E:204:PRO:O	2.19	0.43
1:F:10:ARG:HH21	1:F:14:HIS:CD2	2.36	0.43
1:F:19:LEU:HG	1:F:79:ILE:HG21	2.00	0.43
1:H:105:ASP:O	1:H:109:GLN:HB2	2.18	0.43
1:E:32:TYR:HH	1:E:61:TYR:HH	1.65	0.43
1:E:49:LYS:HE3	1:E:57:PRO:HB2	2.00	0.43
1:E:98:ILE:HD11	1:F:69:ILE:HG23	2.00	0.43
1:E:132:THR:O	1:E:135:GLN:HB3	2.18	0.43
1:G:54:LEU:HD12	1:G:57:PRO:HA	2.01	0.43
1:H:89:THR:OG1	1:H:92:GLU:HB2	2.19	0.43
1:B:206:PRO:HB3	1:B:217:LYS:HB3	2.00	0.43
1:D:56:PHE:O	1:D:70:THR:HG21	2.18	0.43
1:E:141:LEU:O	1:E:144:ARG:HG2	2.18	0.43
1:H:75:ILE:HD12	1:H:75:ILE:H	1.84	0.43
1:H:141:LEU:O	1:H:144:ARG:NH1	2.52	0.43
1:A:10:ARG:CG	1:A:14:HIS:HB2	2.49	0.42
1:C:31:LYS:HE3	1:C:31:LYS:HB2	1.86	0.42
1:D:160:TYR:OH	1:D:197:MET:HE1	2.19	0.42
1:D:180:LEU:O	1:D:184:ILE:HD12	2.18	0.42
1:F:108:ASN:HD22	1:F:108:ASN:HA	1.67	0.42
1:H:144:ARG:HA	1:H:145:PRO:HD3	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:THR:OG1	1:A:92:GLU:HG3	2.19	0.42
1:C:91:GLU:O	1:C:95:ARG:HG3	2.19	0.42
1:E:163:LEU:CB	1:E:184:ILE:HD11	2.48	0.42
1:F:20:LEU:HD12	1:F:79:ILE:HG12	2.01	0.42
1:G:10:ARG:HB3	1:G:206:PRO:O	2.19	0.42
1:G:92:GLU:O	1:G:96:VAL:HG23	2.19	0.42
1:H:82:LYS:O	1:H:83:HIS:CG	2.72	0.42
1:H:140:PHE:HD2	1:H:147:PHE:CD1	2.36	0.42
1:A:216:ASN:HD22	1:A:216:ASN:N	2.17	0.42
1:C:123:LYS:N	1:C:124:PRO:HD2	2.34	0.42
1:E:182:ASP:O	1:E:186:ARG:HB2	2.19	0.42
1:G:2:MET:HE2	1:G:25:SER:HB3	2.01	0.42
1:A:9:ILE:HA	1:A:206:PRO:HG2	2.01	0.42
1:H:130:LEU:HD12	1:H:133:MET:SD	2.59	0.42
1:F:23:THR:HB	1:F:83:HIS:NE2	2.34	0.42
1:G:206:PRO:HA	1:G:217:LYS:NZ	2.34	0.42
1:B:126:TYR:CE1	1:B:133:MET:HE1	2.55	0.42
1:B:138:SER:OG	1:B:180:LEU:HG	2.19	0.42
1:E:144:ARG:HG2	1:E:144:ARG:NH1	2.35	0.42
1:H:43:SER:O	1:H:44:GLN:C	2.62	0.42
1:H:119:PHE:O	1:H:123:LYS:N	2.53	0.42
1:A:99:LEU:HD11	1:A:159:ALA:HB2	2.02	0.42
1:D:16:ILE:HD12	1:D:16:ILE:H	1.84	0.42
1:G:211:VAL:HG22	1:G:211:VAL:O	2.20	0.42
1:H:70:THR:O	1:H:71:GLN:HB2	2.20	0.42
1:H:167:ARG:HH21	1:H:184:ILE:HG21	1.85	0.42
1:D:62:LEU:O	1:D:68:LYS:HA	2.19	0.42
1:D:146:TRP:HB3	1:D:156:ASP:OD2	2.19	0.42
1:F:141:LEU:O	1:F:144:ARG:HG2	2.20	0.42
1:H:42:ARG:O	1:H:46:LEU:HD23	2.19	0.42
1:B:79:ILE:O	1:B:83:HIS:CD2	2.73	0.41
1:C:8:ASP:OD2	1:C:33:THR:HA	2.20	0.41
1:E:33:THR:H	1:E:44:GLN:HE22	1.68	0.41
1:A:125:GLU:O	1:A:128:GLU:HB3	2.19	0.41
1:A:140:PHE:CD2	1:A:147:PHE:HD2	2.32	0.41
1:C:183:PHE:HA	1:C:186:ARG:HH11	1.85	0.41
1:D:21:GLU:CD	1:D:201:ARG:HH22	2.28	0.41
1:D:114:CYS:O	1:D:213:VAL:HG12	2.19	0.41
1:D:144:ARG:HA	1:D:145:PRO:HD3	1.74	0.41
1:E:103:ALA:HB1	1:E:162:VAL:HG21	2.02	0.41
1:G:94:ILE:HD11	1:H:78:TYR:CE1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:130:LEU:HD11	1:H:166:HIS:CD2	2.55	0.41
1:A:63:ILE:HG23	1:A:68:LYS:HD2	2.02	0.41
1:A:92:GLU:O	1:A:96:VAL:HG23	2.20	0.41
1:C:94:ILE:O	1:C:98:ILE:HG13	2.20	0.41
1:G:119:PHE:CZ	1:G:214:TRP:HD1	2.38	0.41
1:G:108:ASN:O	1:G:112:ARG:HB2	2.20	0.41
1:H:114:CYS:HG	1:H:214:TRP:HE3	1.64	0.41
1:B:61:TYR:CD1	1:B:63:ILE:HD11	2.55	0.41
1:B:212:ALA:HB3	1:B:216:ASN:HB3	2.03	0.41
1:F:7:TRP:O	1:F:31:LYS:HG2	2.21	0.41
1:F:146:TRP:O	1:F:149:GLY:N	2.50	0.41
1:F:177:PHE:HA	1:F:178:PRO:HD2	1.84	0.41
1:H:49:LYS:HD3	1:H:57:PRO:HB3	2.02	0.41
1:B:98:ILE:HG22	1:B:99:LEU:N	2.35	0.41
1:C:19:LEU:O	1:C:22:TYR:HB3	2.20	0.41
1:C:21:GLU:HG3	1:C:196:TYR:CG	2.55	0.41
1:E:58:ASN:HB3	1:E:70:THR:CG2	2.50	0.41
1:G:89:THR:OG1	1:G:92:GLU:HG3	2.21	0.41
1:B:208:TYR:HB3	1:B:212:ALA:HB2	2.02	0.41
1:C:70:THR:O	1:C:71:GLN:CB	2.66	0.41
1:C:145:PRO:HG2	1:C:150:ASP:HA	2.03	0.41
1:F:130:LEU:HA	1:F:133:MET:HE3	2.01	0.41
1:B:10:ARG:NE	1:B:165:LEU:HD21	2.36	0.41
1:B:45:TRP:CH2	1:B:61:TYR:HE2	2.38	0.41
1:C:205:LYS:HD3	1:C:217:LYS:HZ3	1.82	0.41
1:D:80:ALA:HB2	1:D:154:PHE:HB3	2.02	0.41
1:E:22:TYR:CD1	1:E:192:LYS:HB2	2.56	0.41
1:H:20:LEU:HD21	1:H:79:ILE:CD1	2.50	0.41
1:A:5:GLY:O	1:A:60:PRO:HA	2.21	0.41
1:B:21:GLU:HG2	1:B:196:TYR:CB	2.49	0.41
1:B:37:ALA:HB1	1:B:38:PRO:CD	2.43	0.41
1:E:20:LEU:HD13	1:E:27:TYR:CD2	2.56	0.41
1:E:43:SER:O	1:E:44:GLN:C	2.63	0.41
1:E:212:ALA:O	1:E:216:ASN:ND2	2.54	0.41
1:F:16:ILE:H	1:F:16:ILE:HG13	1.75	0.41
1:G:4:LEU:HB3	1:G:62:LEU:HD13	2.02	0.41
1:A:69:ILE:HG22	1:A:75:ILE:HG12	2.03	0.41
1:B:206:PRO:HB3	1:B:217:LYS:HB2	2.02	0.41
1:D:71:GLN:O	1:D:75:ILE:HG13	2.21	0.41
1:F:125:GLU:O	1:F:128:GLU:HB2	2.20	0.41
1:F:160:TYR:HD1	1:F:184:ILE:HG22	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:6:TYR:O	1:G:31:LYS:HA	2.21	0.41
1:G:181:LYS:HB3	1:G:181:LYS:HE3	1.89	0.41
1:A:163:LEU:O	1:A:167:ARG:HG2	2.21	0.40
1:A:217:LYS:HE2	1:A:217:LYS:HB2	1.96	0.40
1:C:134:MET:SD	1:C:174:LEU:CD1	3.04	0.40
1:B:5:GLY:O	1:B:60:PRO:HA	2.21	0.40
1:B:7:TRP:CZ3	1:B:34:MET:HE2	2.56	0.40
1:F:2:MET:HE1	1:F:20:LEU:HD11	2.02	0.40
1:A:64:ASP:N	1:A:67:HIS:O	2.53	0.40
1:B:115:TYR:CZ	1:B:208:TYR:HD2	2.39	0.40
1:D:21:GLU:OE2	1:D:196:TYR:HB2	2.21	0.40
1:A:89:THR:HG1	1:A:92:GLU:HG3	1.87	0.40
1:A:95:ARG:NH2	1:A:144:ARG:NH2	2.69	0.40
1:A:130:LEU:N	1:A:131:PRO:HD2	2.36	0.40
1:C:206:PRO:CG	1:C:217:LYS:HB3	2.37	0.40
1:F:10:ARG:HG2	1:F:14:HIS:HB2	2.03	0.40
1:F:96:VAL:O	1:F:100:GLU:HB2	2.22	0.40
1:F:157:PHE:CE2	1:F:187:PHE:HZ	2.39	0.40
1:G:59:LEU:HD23	1:G:59:LEU:HA	1.91	0.40
1:G:98:ILE:HD11	1:H:69:ILE:HG23	2.02	0.40
1:H:17:ARG:HH12	1:H:203:LEU:HD23	1.86	0.40
1:E:108:ASN:O	1:E:112:ARG:HB2	2.22	0.40
1:G:64:ASP:O	1:G:66:ALA:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	215/217 (99%)	184 (86%)	24 (11%)	7 (3%)	3	19
1	B	215/217 (99%)	176 (82%)	28 (13%)	11 (5%)	1	11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	215/217 (99%)	180 (84%)	29 (14%)	6 (3%)	4	21
1	D	215/217 (99%)	183 (85%)	27 (13%)	5 (2%)	5	25
1	E	215/217 (99%)	188 (87%)	20 (9%)	7 (3%)	3	19
1	F	215/217 (99%)	183 (85%)	25 (12%)	7 (3%)	3	19
1	G	215/217 (99%)	182 (85%)	23 (11%)	10 (5%)	2	12
1	H	215/217 (99%)	184 (86%)	22 (10%)	9 (4%)	2	14
All	All	1720/1736 (99%)	1460 (85%)	198 (12%)	62 (4%)	2	17

All (62) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	B	147	PHE
1	B	204	PRO
1	C	14	HIS
1	C	71	GLN
1	E	71	GLN
1	F	119	PHE
1	G	206	PRO
1	A	147	PHE
1	B	40	TYR
1	B	41	ASP
1	B	71	GLN
1	B	142	GLY
1	C	49	LYS
1	D	204	PRO
1	E	65	GLY
1	E	204	PRO
1	F	147	PHE
1	G	142	GLY
1	H	37	ALA
1	B	37	ALA
1	B	184	ILE
1	D	40	TYR
1	E	210	ARG
1	G	41	ASP
1	H	36	ASP
1	H	40	TYR
1	H	188	GLU
1	H	191	GLU

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Mol	Chain	Res	Type
1	H	204	PRO
1	A	57	PRO
1	A	65	GLY
1	B	6	TYR
1	B	119	PHE
1	F	46	LEU
1	G	52	LEU
1	H	71	GLN
1	A	40	TYR
1	A	210	ARG
1	B	72	SER
1	C	37	ALA
1	C	145	PRO
1	D	45	TRP
1	D	141	LEU
1	E	52	LEU
1	G	40	TYR
1	G	215	GLY
1	H	25	SER
1	A	37	ALA
1	D	37	ALA
1	E	37	ALA
1	F	37	ALA
1	F	71	GLN
1	G	37	ALA
1	G	43	SER
1	G	71	GLN
1	G	65	GLY
1	H	38	PRO
1	F	11	GLY
1	C	171	PRO
1	F	204	PRO
1	E	98	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	195/195 (100%)	184 (94%)	11 (6%)	19	47
1	B	195/195 (100%)	181 (93%)	14 (7%)	13	40
1	C	195/195 (100%)	178 (91%)	17 (9%)	9	32
1	D	195/195 (100%)	177 (91%)	18 (9%)	8	30
1	E	195/195 (100%)	177 (91%)	18 (9%)	8	30
1	F	195/195 (100%)	173 (89%)	22 (11%)	5	21
1	G	195/195 (100%)	177 (91%)	18 (9%)	8	30
1	H	195/195 (100%)	178 (91%)	17 (9%)	9	32
All	All	1560/1560 (100%)	1425 (91%)	135 (9%)	9	32

All (135) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	SER
1	A	36	ASP
1	A	68	LYS
1	A	70	THR
1	A	84	ASN
1	A	99	LEU
1	A	135	GLN
1	A	158	LEU
1	A	163	LEU
1	A	165	LEU
1	A	216	ASN
1	B	3	THR
1	B	9	ILE
1	B	39	ASP
1	B	44	GLN
1	B	59	LEU
1	B	69	ILE
1	B	70	THR
1	B	98	ILE
1	B	114	CYS
1	B	132	THR
1	B	138	SER
1	B	144	ARG
1	B	162	VAL
1	B	217	LYS
1	C	3	THR
1	C	34	MET

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Mol	Chain	Res	Type
1	C	39	ASP
1	C	44	GLN
1	C	68	LYS
1	C	72	SER
1	C	77	CYS
1	C	84	ASN
1	C	92	GLU
1	C	120	GLU
1	C	123	LYS
1	C	134	MET
1	C	177	PHE
1	C	179	ASN
1	C	184	ILE
1	C	192	LYS
1	C	205	LYS
1	D	2	MET
1	D	9	ILE
1	D	21	GLU
1	D	24	ASP
1	D	26	SER
1	D	42	ARG
1	D	72	SER
1	D	97	ASP
1	D	98	ILE
1	D	113	VAL
1	D	127	LEU
1	D	132	THR
1	D	134	MET
1	D	167	ARG
1	D	179	ASN
1	D	192	LYS
1	D	201	ARG
1	D	211	VAL
1	E	1	SER
1	E	2	MET
1	E	12	LEU
1	E	41	ASP
1	E	42	ARG
1	E	52	LEU
1	E	84	ASN
1	E	93	LYS
1	E	118	ASP

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Mol	Chain	Res	Type
1	E	136	HIS
1	E	137	PHE
1	E	144	ARG
1	E	179	ASN
1	E	182	ASP
1	E	197	MET
1	E	198	LYS
1	E	205	LYS
1	E	210	ARG
1	F	20	LEU
1	F	26	SER
1	F	47	ASN
1	F	55	ASP
1	F	79	ILE
1	F	84	ASN
1	F	99	LEU
1	F	101	ASN
1	F	104	MET
1	F	105	ASP
1	F	106	VAL
1	F	122	LEU
1	F	123	LYS
1	F	132	THR
1	F	148	VAL
1	F	165	LEU
1	F	182	ASP
1	F	184	ILE
1	F	203	LEU
1	F	205	LYS
1	F	207	LEU
1	F	217	LYS
1	G	4	LEU
1	G	10	ARG
1	G	12	LEU
1	G	16	ILE
1	G	39	ASP
1	G	47	ASN
1	G	84	ASN
1	G	90	GLU
1	G	99	LEU
1	G	112	ARG
1	G	143	LYS

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Mol	Chain	Res	Type
1	G	153	THR
1	G	158	LEU
1	G	165	LEU
1	G	175	ASP
1	G	184	ILE
1	G	207	LEU
1	G	217	LYS
1	H	31	LYS
1	H	33	THR
1	H	44	GLN
1	H	49	LYS
1	H	52	LEU
1	H	61	TYR
1	H	70	THR
1	H	71	GLN
1	H	121	LYS
1	H	127	LEU
1	H	151	LYS
1	H	167	ARG
1	H	175	ASP
1	H	179	ASN
1	H	185	SER
1	H	209	THR
1	H	217	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	HIS
1	A	109	GLN
1	A	135	GLN
1	A	139	GLN
1	A	166	HIS
1	A	179	ASN
1	A	216	ASN
1	B	47	ASN
1	B	83	HIS
1	B	172	ASN
1	C	83	HIS
1	C	102	GLN
1	C	108	ASN
1	C	179	ASN

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Mol	Chain	Res	Type
1	D	14	HIS
1	D	108	ASN
1	D	136	HIS
1	D	179	ASN
1	D	216	ASN
1	E	44	GLN
1	E	71	GLN
1	E	136	HIS
1	E	179	ASN
1	E	216	ASN
1	F	14	HIS
1	F	44	GLN
1	F	108	ASN
1	F	179	ASN
1	G	71	GLN
1	G	108	ASN
1	G	139	GLN
1	H	14	HIS
1	H	83	HIS
1	H	84	ASN
1	H	179	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	217/217 (100%)	0.60	16 (7%) 20 15	6, 18, 29, 35	0
1	B	217/217 (100%)	0.58	8 (3%) 45 31	5, 17, 28, 35	0
1	C	217/217 (100%)	0.58	8 (3%) 45 31	5, 18, 28, 34	0
1	D	217/217 (100%)	0.64	12 (5%) 30 21	5, 17, 28, 35	0
1	E	217/217 (100%)	0.63	13 (5%) 27 19	6, 17, 29, 34	0
1	F	217/217 (100%)	0.57	7 (3%) 50 34	5, 18, 29, 34	0
1	G	217/217 (100%)	0.59	15 (6%) 23 16	5, 17, 29, 34	0
1	H	217/217 (100%)	0.64	10 (4%) 37 25	4, 17, 28, 34	0
All	All	1736/1736 (100%)	0.60	89 (5%) 33 22	4, 18, 29, 35	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	38	PRO	4.6
1	C	124	PRO	3.9
1	D	116	SER	3.9
1	F	182	ASP	3.7
1	E	131	PRO	3.6
1	A	136	HIS	3.6
1	A	216	ASN	3.5
1	G	131	PRO	3.4
1	D	75	ILE	3.4
1	B	117	PRO	3.3
1	G	103	ALA	3.3
1	F	206	PRO	3.2
1	G	199	SER	3.2
1	D	96	VAL	3.2
1	E	212	ALA	3.2
1	E	132	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	32	TYR	3.1
1	A	37	ALA	3.0
1	A	40	TYR	3.0
1	H	1	SER	3.0
1	B	37	ALA	3.0
1	C	80	ALA	3.0
1	A	105	ASP	2.9
1	E	133	MET	2.9
1	E	185	SER	2.8
1	A	33	THR	2.8
1	D	184	ILE	2.8
1	G	86	CYS	2.8
1	A	111	ALA	2.8
1	H	92	GLU	2.8
1	G	189	GLY	2.8
1	E	108	ASN	2.8
1	F	1	SER	2.8
1	C	119	PHE	2.7
1	H	161	ASP	2.7
1	F	170	GLU	2.7
1	E	80	ALA	2.7
1	A	215	GLY	2.7
1	A	80	ALA	2.7
1	G	132	THR	2.6
1	B	38	PRO	2.6
1	D	35	GLY	2.6
1	F	111	ALA	2.6
1	A	27	TYR	2.5
1	F	131	PRO	2.5
1	C	142	GLY	2.5
1	D	118	ASP	2.5
1	C	35	GLY	2.4
1	A	159	ALA	2.4
1	H	111	ALA	2.4
1	A	77	CYS	2.4
1	D	215	GLY	2.4
1	C	37	ALA	2.4
1	B	63	ILE	2.3
1	G	27	TYR	2.3
1	D	145	PRO	2.3
1	A	172	ASN	2.3
1	E	206	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	103	ALA	2.3
1	G	118	ASP	2.3
1	C	178	PRO	2.3
1	D	101	ASN	2.3
1	H	58	ASN	2.3
1	H	38	PRO	2.2
1	H	206	PRO	2.2
1	B	132	THR	2.2
1	D	3	THR	2.2
1	G	128	GLU	2.2
1	A	30	LYS	2.2
1	H	23	THR	2.2
1	G	216	ASN	2.2
1	B	5	GLY	2.2
1	E	35	GLY	2.2
1	F	156	ASP	2.1
1	G	177	PHE	2.1
1	B	136	HIS	2.1
1	B	138	SER	2.1
1	E	177	PHE	2.1
1	D	84	ASN	2.1
1	E	136	HIS	2.1
1	G	24	ASP	2.1
1	E	195	ALA	2.1
1	G	71	GLN	2.1
1	G	178	PRO	2.1
1	E	171	PRO	2.0
1	G	153	THR	2.0
1	H	37	ALA	2.0
1	H	124	PRO	2.0
1	A	46	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.