



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:43 AM UTC

PDB ID : 1KNX / pdb_00001knx
Title : HPr kinase/phosphatase from Mycoplasma pneumoniae
Authors : Allen, G.S.
Deposited on : 2001-12-19
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

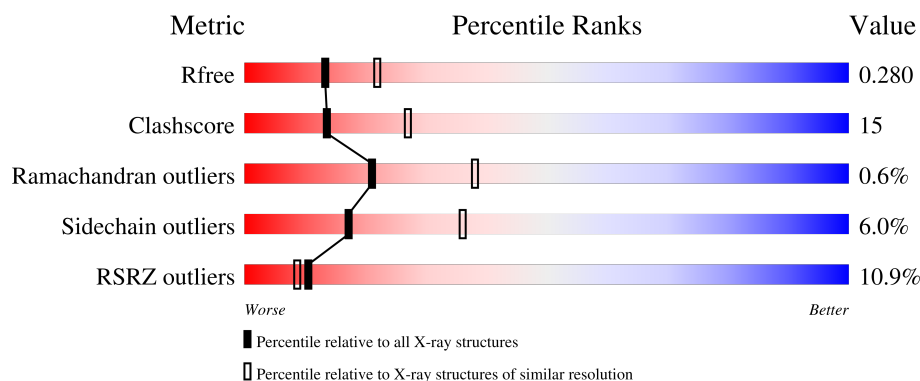
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	312	25% 56% 36% 5%
1	B	312	2% 65% 30%
1	C	312	4% 64% 29%
1	D	312	23% 64% 30%
1	E	312	3% 70% 25%

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Mol	Chain	Length	Quality of chain
1	F	312	7% 68% 28% • •

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 14717 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 Probable HPr(Ser) kinase/phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	0	0	0
			2392	1546	404	436	6			
1	B	304	Total	C	N	O	S	0	0	0
			2415	1559	409	441	6			
1	C	300	Total	C	N	O	S	0	0	0
			2375	1532	403	434	6			
1	D	304	Total	C	N	O	S	0	0	0
			2400	1548	408	438	6			
1	E	310	Total	C	N	O	S	0	0	0
			2446	1576	414	450	6			
1	F	306	Total	C	N	O	S	0	0	0
			2407	1555	404	442	6			

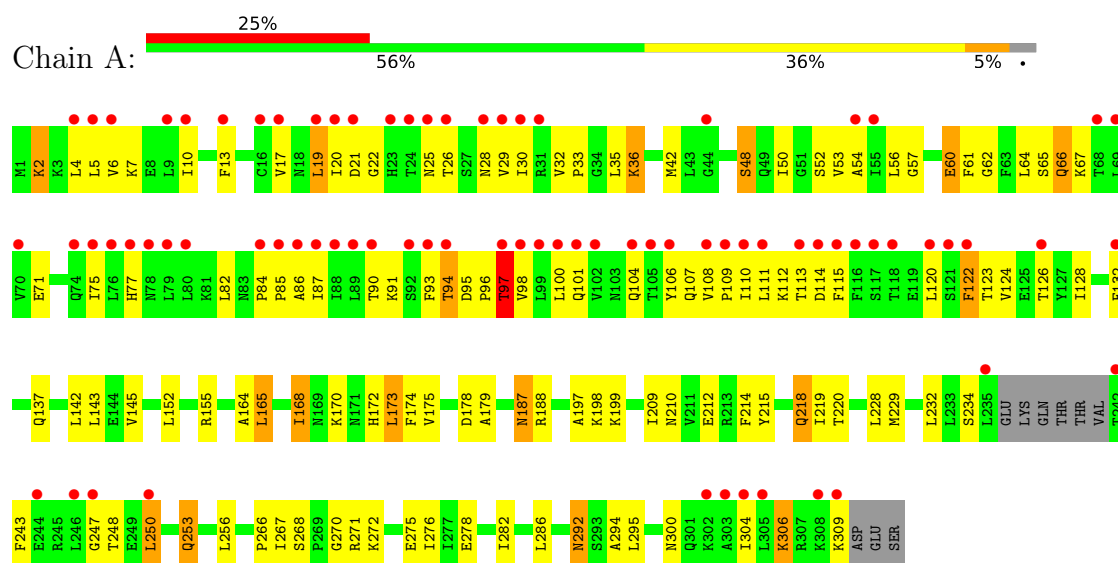
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	34	Total	O	0	0
			34	34		
2	B	55	Total	O	0	0
			55	55		
2	C	38	Total	O	0	0
			38	38		
2	D	39	Total	O	0	0
			39	39		
2	E	61	Total	O	0	0
			61	61		
2	F	55	Total	O	0	0
			55	55		

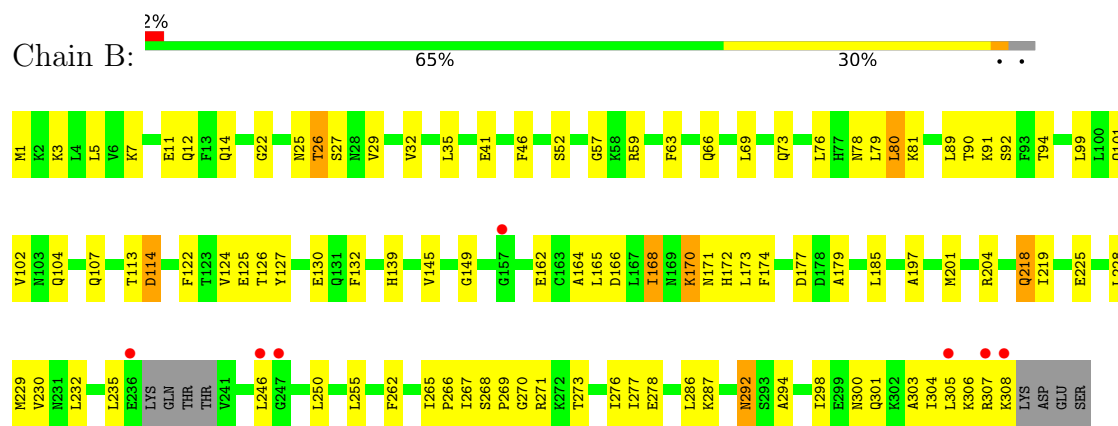
3 Residue-property plots [i](#)

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.

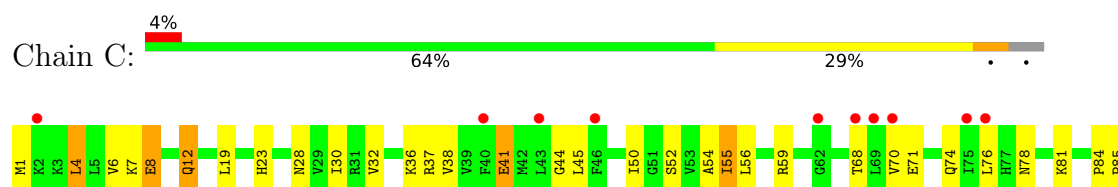
- Molecule 1: Probable HPr(Ser) kinase/phosphatase

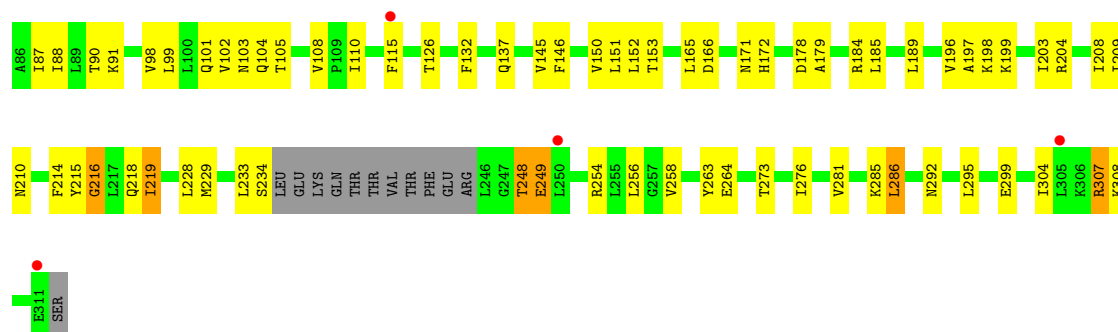


- Molecule 1: Probable HPr(Ser) kinase/phosphatase

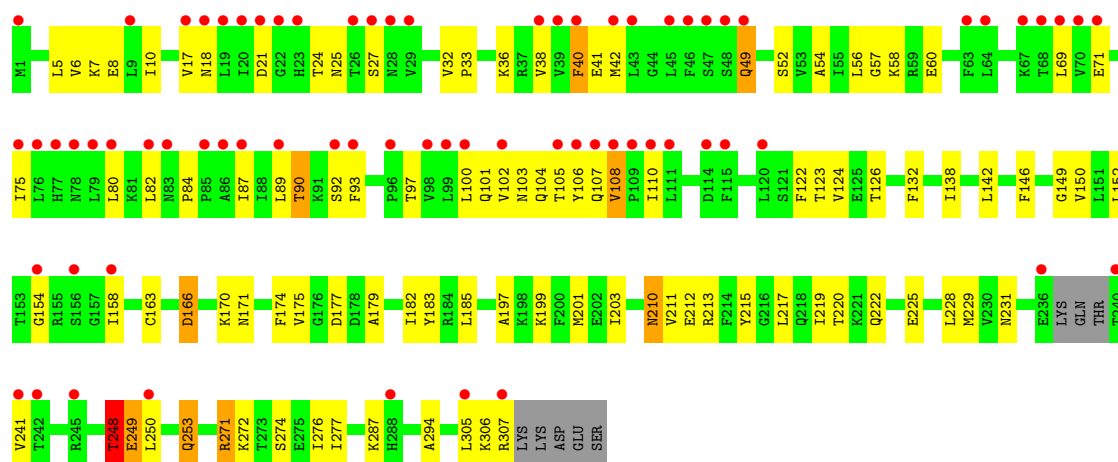


- Molecule 1: Probable HPr(Ser) kinase/phosphatase

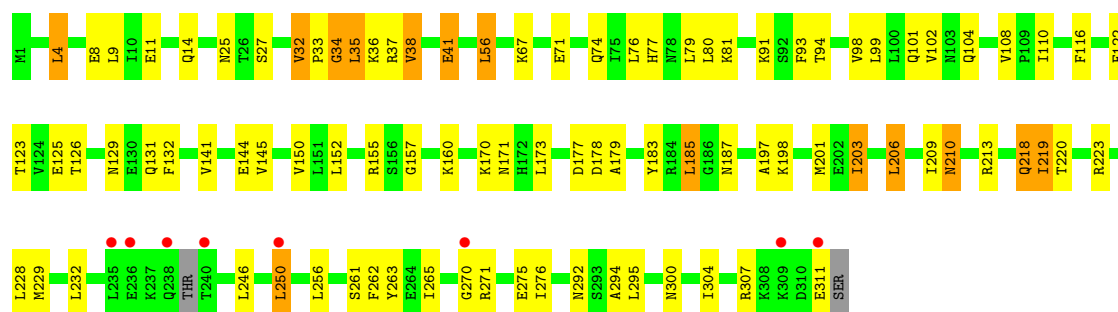




- Molecule 1: Probable HPr(Ser) kinase/phosphatase

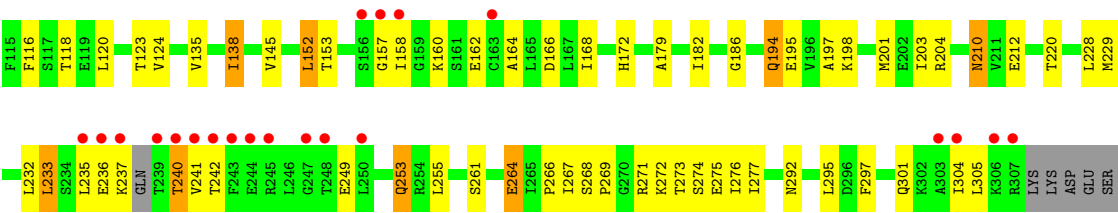


- Molecule 1: Probable HPr(Ser) kinase/phosphatase



- Molecule 1: Probable HPr(Ser) kinase/phosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	117.73Å 127.84Å 170.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.22 – 2.50 29.22 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.3 (29.22-2.50) 91.2 (29.22-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.280 0.229 , 0.280	Depositor DCC
R_{free} test set	4908 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	41.2	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14717	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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	0.42	0/2427	0.82	2/3270 (0.1%)
1	B	0.48	0/2450	0.89	4/3301 (0.1%)
1	C	0.44	0/2409	0.88	3/3245 (0.1%)
1	D	0.45	0/2435	0.89	6/3283 (0.2%)
1	E	0.53	0/2481	0.91	8/3346 (0.2%)
1	F	0.56	2/2442 (0.1%)	0.94	8/3295 (0.2%)
All	All	0.48	2/14644 (0.0%)	0.89	31/19740 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	F	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	233	LEU	C-N	-10.11	1.19	1.33
1	F	237	LYS	N-CA	5.09	1.55	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	198	LYS	CB-CA-C	-7.41	108.02	116.54
1	D	71	GLU	N-CA-C	-7.19	103.05	111.03
1	F	235	LEU	N-CA-C	6.95	119.70	111.71
1	F	94	THR	N-CA-C	6.82	118.37	111.07
1	F	242	THR	N-CA-C	-6.54	99.55	109.95
1	B	170	LYS	N-CA-C	-6.42	105.19	113.16
1	F	233	LEU	CA-C-N	6.38	132.39	121.86
1	F	233	LEU	C-N-CA	6.38	132.39	121.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	240	THR	N-CA-C	-6.24	105.98	113.97
1	D	108	VAL	N-CA-C	6.17	113.66	107.55
1	E	36	LYS	N-CA-C	6.01	119.06	110.10
1	D	241	VAL	N-CA-C	5.99	117.80	109.29
1	B	94	THR	N-CA-C	5.98	120.19	113.01
1	B	124	VAL	N-CA-C	5.89	116.08	110.42
1	E	219	ILE	N-CA-C	5.65	117.87	111.88
1	C	216	GLY	N-CA-C	5.65	119.80	112.68
1	C	150	VAL	N-CA-C	5.65	116.00	107.75
1	E	32	VAL	CA-C-N	5.62	125.60	120.21
1	E	32	VAL	C-N-CA	5.62	125.60	120.21
1	E	187	ASN	N-CA-C	-5.54	105.83	112.59
1	C	307	ARG	N-CA-C	-5.50	106.58	113.72
1	B	80	LEU	N-CA-C	5.43	117.95	111.71
1	A	170	LYS	N-CA-C	-5.41	105.99	112.59
1	F	194	GLN	N-CA-C	-5.32	103.36	110.55
1	E	94	THR	N-CA-C	5.31	119.47	112.89
1	D	36	LYS	N-CA-C	5.27	118.20	110.30
1	E	34	GLY	N-CA-C	5.25	118.83	112.48
1	D	248	THR	N-CA-C	-5.11	106.21	112.90
1	D	276	ILE	N-CA-C	-5.07	105.60	110.72
1	F	36	LYS	N-CA-C	5.06	117.64	110.10
1	A	36	LYS	N-CA-C	5.03	117.53	110.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	236	GLU	Peptide,Mainchain

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	2392	0	2488	111	0
1	B	2415	0	2520	67	0
1	C	2375	0	2478	73	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2400	0	2490	75	0
1	E	2446	0	2528	61	0
1	F	2407	0	2491	64	0
2	A	34	0	0	0	0
2	B	55	0	0	3	0
2	C	38	0	0	3	0
2	D	39	0	0	0	0
2	E	61	0	0	2	0
2	F	55	0	0	1	0
All	All	14717	0	14995	431	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 15.

All (431) 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:A:282:ILE:HG12	1:B:218:GLN:HG2	1.21	1.11
1:E:292:ASN:HD22	1:E:295:LEU:H	1.05	0.98
1:B:90:THR:HG23	1:B:92:SER:H	1.28	0.97
1:E:292:ASN:ND2	1:E:295:LEU:H	1.64	0.95
1:A:218:GLN:NE2	1:A:218:GLN:H	1.62	0.95
1:A:218:GLN:H	1:A:218:GLN:HE21	0.94	0.93
1:C:78:ASN:HA	1:C:81:LYS:HE2	1.51	0.93
1:A:218:GLN:HE21	1:A:218:GLN:N	1.69	0.90
1:C:292:ASN:HD22	1:C:295:LEU:H	1.21	0.89
1:B:168:ILE:HD11	1:B:174:PHE:HB2	1.55	0.88
1:A:187:ASN:HD22	1:A:187:ASN:H	1.22	0.86
1:C:32:VAL:HB	1:C:52:SER:HB3	1.56	0.86
1:F:49:GLN:HE22	1:F:83:ASN:HD22	1.21	0.84
1:A:250:LEU:HD23	1:A:250:LEU:H	1.41	0.84
1:A:113:THR:HG22	1:A:115:PHE:H	1.44	0.83
1:A:282:ILE:HG12	1:B:218:GLN:CG	2.08	0.83
1:E:108:VAL:O	1:E:110:ILE:HD12	1.79	0.83
1:E:210:ASN:C	1:E:210:ASN:HD22	1.88	0.82
1:F:103:ASN:HD21	1:F:108:VAL:H	1.24	0.82
1:E:218:GLN:H	1:E:218:GLN:HE21	1.29	0.81
1:F:292:ASN:HD22	1:F:295:LEU:H	1.30	0.80
1:A:25:ASN:HB2	1:A:109:PRO:HD3	1.64	0.79
1:F:166:ASP:HB3	1:F:274:SER:HB3	1.63	0.79
1:B:273:THR:O	1:B:276:ILE:HG22	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:LEU:O	1:B:113:THR:HG22	1.84	0.78
1:B:22:GLY:HA2	1:B:107:GLN:HG3	1.67	0.77
1:A:19:LEU:H	1:A:19:LEU:HD23	1.48	0.76
1:D:179:ALA:HB3	1:D:197:ALA:HB2	1.68	0.75
1:A:187:ASN:H	1:A:187:ASN:ND2	1.85	0.74
1:C:55:ILE:HD13	1:C:88:ILE:HG23	1.70	0.74
1:B:1:MET:HG2	1:B:3:LYS:H	1.54	0.72
1:B:35:LEU:HG	1:B:125:GLU:HG2	1.69	0.72
1:F:157:GLY:HA2	1:F:160:LYS:NZ	2.04	0.72
1:D:271:ARG:HG2	1:D:272:LYS:N	2.04	0.72
1:F:273:THR:O	1:F:276:ILE:HG22	1.89	0.72
1:D:183:TYR:HD2	1:D:185:LEU:HD13	1.55	0.71
1:A:126:THR:HG23	1:B:122:PHE:CE2	2.25	0.71
1:D:201:MET:HE3	1:D:203:ILE:HD11	1.72	0.71
1:A:35:LEU:HD13	1:A:120:LEU:HD23	1.73	0.70
1:A:97:THR:HG23	1:A:98:VAL:H	1.56	0.70
1:F:212:GLU:HG3	1:F:220:THR:HG21	1.73	0.69
1:A:271:ARG:NH2	1:A:276:ILE:HG13	2.08	0.68
1:B:76:LEU:HD13	1:B:99:LEU:HA	1.76	0.68
1:B:162:GLU:HG2	1:E:271:ARG:HD2	1.77	0.67
1:B:90:THR:HG23	1:B:92:SER:N	2.06	0.67
1:C:90:THR:HG21	2:C:313:HOH:O	1.96	0.66
1:A:56:LEU:HA	1:A:60:GLU:HG3	1.78	0.66
1:D:123:THR:HG22	1:D:124:VAL:N	2.10	0.66
1:A:126:THR:HG23	1:B:122:PHE:HE2	1.60	0.65
1:B:270:GLY:HA2	1:D:271:ARG:O	1.96	0.65
1:B:246:LEU:HD22	1:B:298:ILE:HD11	1.78	0.65
1:D:56:LEU:HD23	1:D:60:GLU:HG2	1.79	0.65
1:C:90:THR:HG22	1:C:91:LYS:N	2.12	0.65
1:B:292:ASN:C	1:B:292:ASN:HD22	2.05	0.64
1:C:233:LEU:HD11	1:C:264:GLU:OE2	1.97	0.64
1:F:77:HIS:CE1	1:F:102:VAL:HG13	2.32	0.64
1:A:57:GLY:HA2	1:A:90:THR:HG22	1.79	0.64
1:C:292:ASN:ND2	1:C:295:LEU:H	1.94	0.64
1:C:55:ILE:HD12	1:C:56:LEU:H	1.63	0.64
1:D:174:PHE:CD2	1:D:211:VAL:HG21	2.32	0.64
1:C:165:LEU:HD22	1:C:203:ILE:HD13	1.80	0.64
1:B:32:VAL:HB	1:B:52:SER:HB3	1.79	0.63
1:A:36:LYS:HE3	1:B:59:ARG:HD2	1.80	0.63
1:D:32:VAL:HB	1:D:52:SER:HB3	1.80	0.63
1:B:301:GLN:O	1:B:305:LEU:HD13	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:300:ASN:O	1:B:304:ILE:HG12	2.00	0.62
1:B:170:LYS:O	1:B:171:ASN:HB2	1.99	0.61
1:C:198:LYS:O	1:C:199:LYS:HD2	1.99	0.61
1:D:228:LEU:HD23	1:D:229:MET:N	2.15	0.61
1:C:171:ASN:HA	1:C:219:ILE:HD11	1.81	0.61
1:E:292:ASN:HD22	1:E:295:LEU:N	1.88	0.61
1:C:228:LEU:HD11	1:C:263:TYR:CE1	2.36	0.61
1:D:80:LEU:HD22	1:D:108:VAL:HG21	1.83	0.60
1:E:219:ILE:HD12	1:E:220:THR:HG23	1.82	0.60
1:A:110:ILE:C	1:A:111:LEU:HD12	2.26	0.60
1:D:38:VAL:HG22	1:D:41:GLU:HB2	1.82	0.60
1:D:305:LEU:C	1:D:307:ARG:H	2.09	0.60
1:C:145:VAL:HG13	1:C:172:HIS:CD2	2.37	0.60
1:E:35:LEU:HD22	1:E:125:GLU:HB2	1.82	0.60
1:F:5:LEU:HD23	1:F:7:LYS:HE2	1.83	0.60
1:A:32:VAL:HB	1:A:52:SER:HB3	1.82	0.60
1:A:232:LEU:HB3	1:A:267:ILE:HD11	1.83	0.60
1:B:7:LYS:HB3	1:B:26:THR:HG23	1.84	0.59
1:F:210:ASN:C	1:F:210:ASN:HD22	2.10	0.59
1:A:17:VAL:HB	1:A:111:LEU:HD23	1.85	0.59
1:B:12:GLN:HG3	1:B:127:TYR:CE1	2.38	0.59
1:F:201:MET:HE3	1:F:203:ILE:HD11	1.84	0.59
1:F:153:THR:HG21	1:F:229:MET:HE2	1.85	0.58
1:E:179:ALA:HB3	1:E:197:ALA:HB2	1.85	0.58
1:C:210:ASN:HD22	1:C:210:ASN:C	2.11	0.58
1:A:179:ALA:HB3	1:A:197:ALA:HB2	1.85	0.58
1:A:67:LYS:HG3	1:A:71:GLU:OE2	2.03	0.58
1:A:22:GLY:HA2	1:A:107:GLN:HG3	1.85	0.58
1:F:52:SER:O	1:F:85:PRO:HD2	2.04	0.58
1:A:292:ASN:HD21	1:A:294:ALA:HB3	1.69	0.57
1:B:177:ASP:OD1	1:B:201:MET:HG3	2.04	0.57
1:A:61:PHE:HB2	1:A:93:PHE:HD1	1.69	0.57
1:A:209:ILE:HG13	1:A:214:PHE:CZ	2.39	0.57
1:B:179:ALA:HB3	1:B:197:ALA:HB2	1.86	0.57
1:C:228:LEU:HD11	1:C:263:TYR:CD1	2.40	0.57
1:E:37:ARG:HD3	2:E:342:HOH:O	2.04	0.57
1:E:157:GLY:HA2	1:E:160:LYS:HE3	1.85	0.57
1:F:138:ILE:HD11	1:F:182:ILE:HG23	1.86	0.57
1:F:158:ILE:HG21	1:F:204:ARG:HG3	1.85	0.57
1:C:101:GLN:O	1:C:104:GLN:HG2	2.05	0.56
1:E:34:GLY:HA2	1:E:129:ASN:HD21	1.69	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:GLU:CD	1:C:41:GLU:H	2.12	0.56
1:D:57:GLY:HA2	1:D:90:THR:OG1	2.04	0.56
1:D:150:VAL:HG12	1:D:152:LEU:HD12	1.86	0.56
1:F:179:ALA:O	1:F:194:GLN:HG3	2.06	0.56
1:E:292:ASN:HD21	1:E:294:ALA:HB3	1.70	0.56
1:C:90:THR:HG22	1:C:91:LYS:H	1.70	0.56
1:A:93:PHE:O	1:A:94:THR:HG22	2.06	0.56
1:A:292:ASN:ND2	1:A:295:LEU:H	2.04	0.56
1:D:6:VAL:O	1:D:10:ILE:HG13	2.06	0.56
1:A:54:ALA:HB3	1:A:87:ILE:HD13	1.88	0.56
1:C:248:THR:HG22	1:C:295:LEU:HG	1.87	0.55
1:E:228:LEU:HD12	1:E:261:SER:HB2	1.87	0.55
1:E:201:MET:HE3	1:E:203:ILE:CD1	2.37	0.55
1:B:69:LEU:O	1:B:73:GLN:HG3	2.06	0.55
1:E:210:ASN:C	1:E:210:ASN:ND2	2.59	0.55
1:D:199:LYS:O	1:D:211:VAL:HG22	2.07	0.55
1:A:152:LEU:HG	1:A:232:LEU:HD21	1.89	0.55
1:D:210:ASN:HD21	1:D:212:GLU:HB3	1.72	0.55
1:A:172:HIS:HE1	1:A:278:GLU:OE2	1.89	0.54
1:C:55:ILE:HD12	1:C:56:LEU:N	2.22	0.54
1:A:271:ARG:CZ	1:A:276:ILE:HG13	2.36	0.54
1:B:168:ILE:CD1	1:B:174:PHE:HB2	2.31	0.54
1:C:44:GLY:C	1:C:45:LEU:HD12	2.33	0.54
1:B:145:VAL:HG13	1:B:172:HIS:CD2	2.42	0.54
1:D:249:GLU:HG3	1:D:250:LEU:H	1.71	0.54
1:A:33:PRO:HB3	1:A:132:PHE:CD1	2.43	0.54
1:D:84:PRO:HG3	1:D:87:ILE:HD11	1.89	0.54
1:E:218:GLN:HE21	1:E:218:GLN:N	2.02	0.54
1:D:42:MET:O	1:D:75:ILE:HG23	2.07	0.54
1:A:143:LEU:HD21	1:A:164:ALA:HB1	1.90	0.54
1:A:268:SER:HB3	1:F:204:ARG:HH22	1.73	0.54
1:E:67:LYS:HB3	1:E:71:GLU:HB2	1.89	0.54
1:E:206:LEU:HD23	1:E:209:ILE:HG21	1.90	0.54
1:B:265:ILE:HG23	1:B:266:PRO:HD2	1.90	0.53
1:D:110:ILE:HG13	1:D:110:ILE:O	2.09	0.53
1:E:250:LEU:HD12	1:E:250:LEU:H	1.72	0.53
1:E:304:ILE:O	1:E:307:ARG:HG2	2.08	0.53
1:C:249:GLU:CD	1:C:249:GLU:H	2.16	0.53
1:D:166:ASP:HB3	1:D:274:SER:HB3	1.90	0.53
1:A:187:ASN:HD22	1:A:187:ASN:N	1.90	0.53
1:B:294:ALA:O	1:B:298:ILE:HG12	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:MET:HE1	1:C:132:PHE:HA	1.90	0.53
1:B:41:GLU:HG3	1:B:46:PHE:HB2	1.91	0.52
1:F:233:LEU:HD11	1:F:264:GLU:HG2	1.92	0.52
1:A:7:LYS:HG3	1:A:26:THR:O	2.09	0.52
1:D:38:VAL:HG22	1:D:38:VAL:O	2.09	0.52
1:A:145:VAL:HG13	1:A:172:HIS:CD2	2.44	0.52
1:C:153:THR:HG21	1:C:229:MET:HE2	1.91	0.52
1:C:36:LYS:HB2	1:C:52:SER:OG	2.09	0.52
1:D:183:TYR:CD2	1:D:185:LEU:HD13	2.38	0.52
1:B:139:HIS:HB3	2:B:353:HOH:O	2.09	0.52
1:C:68:THR:HG22	1:C:70:VAL:H	1.74	0.52
1:B:1:MET:HE1	1:B:132:PHE:CE1	2.45	0.52
1:A:56:LEU:HD13	1:A:60:GLU:HG3	1.90	0.52
1:C:54:ALA:O	1:C:87:ILE:HA	2.10	0.52
1:D:33:PRO:HB3	1:D:132:PHE:CD1	2.46	0.51
1:C:19:LEU:HD21	1:C:23:HIS:HA	1.91	0.51
1:A:168:ILE:HG13	1:A:219:ILE:HB	1.93	0.51
1:C:76:LEU:HD13	1:C:99:LEU:HA	1.92	0.51
1:D:54:ALA:HB3	1:D:87:ILE:HG12	1.92	0.51
1:F:164:ALA:O	1:F:168:ILE:HG13	2.10	0.51
1:A:168:ILE:CG1	1:A:219:ILE:HD13	2.41	0.51
1:B:130:GLU:HG2	1:B:185:LEU:HD11	1.93	0.51
1:F:32:VAL:HB	1:F:52:SER:HB3	1.93	0.51
1:B:91:LYS:HE2	1:B:114:ASP:OD1	2.11	0.51
1:C:98:VAL:O	1:C:102:VAL:HG23	2.11	0.51
1:E:76:LEU:HD13	1:E:99:LEU:HA	1.92	0.51
1:A:152:LEU:HG	1:A:232:LEU:CD2	2.41	0.50
1:D:41:GLU:HG2	1:D:82:LEU:CD1	2.41	0.50
1:D:248:THR:HG22	1:D:294:ALA:HB1	1.93	0.50
1:E:101:GLN:O	1:E:104:GLN:HG3	2.11	0.50
1:F:301:GLN:O	1:F:305:LEU:HG	2.12	0.50
1:F:123:THR:OG1	1:F:124:VAL:N	2.43	0.50
1:E:34:GLY:H	1:E:129:ASN:ND2	2.10	0.50
1:A:247:GLY:H	1:A:294:ALA:HB2	1.76	0.50
1:A:60:GLU:CD	1:A:60:GLU:H	2.20	0.49
1:D:248:THR:HG22	1:D:294:ALA:CB	2.42	0.49
1:A:188:ARG:HG2	1:A:188:ARG:HH11	1.77	0.49
1:A:292:ASN:HD22	1:A:295:LEU:H	1.59	0.49
1:B:218:GLN:H	1:B:218:GLN:NE2	2.10	0.49
1:A:96:PRO:O	1:A:100:LEU:HG	2.11	0.49
1:B:3:LYS:HE3	1:B:29:VAL:HG11	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:69:LEU:N	1:D:69:LEU:HD22	2.27	0.49
1:E:228:LEU:HD11	1:E:263:TYR:CE2	2.47	0.49
1:E:229:MET:HE2	1:E:262:PHE:CD1	2.48	0.49
1:A:7:LYS:HG2	1:A:26:THR:HB	1.94	0.49
1:A:270:GLY:HA2	1:F:271:ARG:O	2.13	0.49
1:C:295:LEU:O	1:C:299:GLU:HG2	2.13	0.49
1:D:210:ASN:C	1:D:210:ASN:HD22	2.21	0.49
1:D:219:ILE:HD12	1:D:220:THR:HG23	1.94	0.49
1:A:56:LEU:HD23	1:A:87:ILE:HG23	1.95	0.49
1:E:11:GLU:O	1:E:14:GLN:HG3	2.12	0.48
1:F:49:GLN:NE2	1:F:83:ASN:HD22	2.02	0.48
1:F:228:LEU:HD12	1:F:261:SER:HB2	1.94	0.48
1:F:292:ASN:ND2	1:F:295:LEU:HG	2.28	0.48
1:A:4:LEU:HG	1:A:30:ILE:HD12	1.95	0.48
1:E:201:MET:HE3	1:E:203:ILE:HD13	1.95	0.48
1:A:228:LEU:HD23	1:A:229:MET:N	2.28	0.48
1:A:187:ASN:ND2	1:A:187:ASN:N	2.48	0.48
1:B:271:ARG:HG2	1:D:166:ASP:OD1	2.13	0.48
1:C:204:ARG:NH2	1:F:268:SER:HB2	2.28	0.48
1:A:42:MET:HB3	1:A:75:ILE:CG2	2.44	0.48
1:C:103:ASN:C	1:C:105:THR:H	2.22	0.48
1:F:89:LEU:O	1:F:113:THR:HG22	2.14	0.48
1:F:91:LYS:HB3	1:F:114:ASP:HA	1.96	0.48
1:B:228:LEU:CD1	1:B:230:VAL:HG23	2.44	0.47
1:D:177:ASP:OD1	1:D:201:MET:HG3	2.14	0.47
1:D:253:GLN:HE21	1:D:253:GLN:C	2.22	0.47
1:A:57:GLY:HA2	1:A:90:THR:CG2	2.43	0.47
1:C:68:THR:HB	1:C:71:GLU:HG3	1.96	0.47
1:F:77:HIS:HD2	1:F:106:TYR:OH	1.97	0.47
1:A:48:SER:O	1:A:82:LEU:HD13	2.14	0.47
1:D:102:VAL:O	1:D:106:TYR:HD1	1.97	0.47
1:B:303:ALA:O	1:B:307:ARG:HG3	2.14	0.47
1:C:90:THR:CG2	1:C:91:LYS:N	2.77	0.47
1:F:138:ILE:CD1	1:F:182:ILE:HG23	2.45	0.47
1:F:157:GLY:HA2	1:F:160:LYS:HZ3	1.75	0.47
1:A:271:ARG:HD2	1:F:162:GLU:HB3	1.95	0.47
1:B:11:GLU:O	1:B:14:GLN:HG3	2.14	0.47
1:B:25:ASN:C	1:B:27:SER:H	2.23	0.47
1:B:78:ASN:HA	1:B:81:LYS:HE2	1.95	0.47
1:C:108:VAL:O	1:C:110:ILE:HG13	2.14	0.47
1:C:90:THR:HG23	1:C:115:PHE:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:91:LYS:HD3	1:F:116:PHE:CZ	2.50	0.47
1:A:95:ASP:C	1:A:97:THR:H	2.23	0.47
1:D:7:LYS:HA	1:D:10:ILE:HD12	1.97	0.47
1:D:69:LEU:HD22	1:D:69:LEU:H	1.80	0.47
1:E:229:MET:HE2	1:E:262:PHE:CE1	2.50	0.47
1:A:168:ILE:HG13	1:A:219:ILE:HD13	1.96	0.47
1:C:4:LEU:HD22	1:C:30:ILE:HD12	1.97	0.47
1:C:184:ARG:O	1:C:185:LEU:HD12	2.15	0.47
1:D:25:ASN:C	1:D:27:SER:N	2.72	0.46
1:E:37:ARG:HA	2:E:316:HOH:O	2.15	0.46
1:A:304:ILE:C	1:A:306:LYS:H	2.23	0.46
1:B:101:GLN:O	1:B:104:GLN:HB2	2.15	0.46
1:C:204:ARG:HH22	1:F:268:SER:HB2	1.80	0.46
1:D:60:GLU:OE2	1:D:60:GLU:N	2.48	0.46
1:A:32:VAL:CB	1:A:52:SER:HB3	2.46	0.46
1:A:91:LYS:C	1:A:93:PHE:H	2.23	0.46
1:B:168:ILE:HG23	1:B:219:ILE:HG12	1.98	0.46
1:E:250:LEU:HD12	1:E:250:LEU:N	2.31	0.46
1:C:281:VAL:O	1:C:285:LYS:HG3	2.16	0.46
1:E:183:TYR:CD1	1:E:185:LEU:HD13	2.51	0.46
1:E:218:GLN:H	1:E:218:GLN:NE2	2.06	0.46
1:F:195:GLU:O	1:F:198:LYS:HG3	2.15	0.46
1:A:2:LYS:HE3	1:A:2:LYS:HA	1.97	0.46
1:C:254:ARG:HA	1:C:258:VAL:O	2.16	0.46
1:D:138:ILE:HD13	1:D:182:ILE:HD13	1.97	0.46
1:D:7:LYS:HG3	1:D:8:GLU:H	1.81	0.46
1:E:77:HIS:O	1:E:81:LYS:HB2	2.15	0.46
1:D:215:TYR:HB2	1:D:219:ILE:HG12	1.98	0.46
1:E:32:VAL:HG13	1:E:33:PRO:HD2	1.98	0.46
1:A:122:PHE:HD1	1:A:123:THR:HG23	1.80	0.46
1:F:272:LYS:HD2	1:F:275:GLU:OE2	2.15	0.46
1:B:80:LEU:HD12	1:B:102:VAL:HG12	1.97	0.45
1:C:50:ILE:O	1:C:85:PRO:HD3	2.16	0.45
1:C:179:ALA:HB3	1:C:197:ALA:HB2	1.98	0.45
1:A:33:PRO:HD3	1:A:132:PHE:CD2	2.52	0.45
1:B:232:LEU:HD21	1:B:277:ILE:HD11	1.98	0.45
1:A:101:GLN:HA	1:A:104:GLN:HE21	1.82	0.45
1:E:232:LEU:HA	1:E:265:ILE:O	2.16	0.45
1:C:273:THR:HA	1:C:276:ILE:HD12	1.98	0.45
1:E:219:ILE:HD12	1:E:219:ILE:C	2.42	0.45
1:C:59:ARG:NH2	1:D:40:PHE:HB2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:209:ILE:O	1:C:209:ILE:HG13	2.17	0.45
1:C:218:GLN:HB2	1:D:146:PHE:CD2	2.51	0.45
1:F:25:ASN:C	1:F:27:SER:H	2.22	0.45
1:F:77:HIS:ND1	1:F:102:VAL:HG13	2.32	0.45
1:A:243:PHE:CE2	1:A:266:PRO:HD3	2.52	0.45
1:D:170:LYS:O	1:D:171:ASN:HB2	2.15	0.45
1:F:118:THR:HG23	2:F:322:HOH:O	2.16	0.45
1:B:229:MET:HE2	1:B:262:PHE:HE1	1.82	0.45
1:C:55:ILE:CD1	1:C:88:ILE:HG23	2.44	0.45
1:D:25:ASN:C	1:D:27:SER:H	2.24	0.45
1:D:69:LEU:H	1:D:69:LEU:CD2	2.30	0.45
1:A:10:ILE:HD13	1:A:111:LEU:HD21	1.99	0.45
1:A:19:LEU:H	1:A:19:LEU:CD2	2.24	0.45
1:B:41:GLU:HG3	1:B:46:PHE:CB	2.47	0.45
1:C:214:PHE:HB3	1:C:215:TYR:CE1	2.52	0.45
1:D:17:VAL:HG12	1:D:18:ASN:N	2.32	0.45
1:D:103:ASN:C	1:D:105:THR:N	2.75	0.45
1:D:305:LEU:C	1:D:307:ARG:N	2.75	0.45
1:A:10:ILE:CD1	1:A:111:LEU:HD21	2.47	0.44
1:B:300:ASN:HB3	2:B:336:HOH:O	2.18	0.44
1:C:145:VAL:O	1:C:146:PHE:HB2	2.17	0.44
1:D:38:VAL:CG2	1:D:41:GLU:HB2	2.47	0.44
1:A:62:GLY:HA2	1:A:65:SER:HB3	1.98	0.44
1:A:106:TYR:HB3	1:A:108:VAL:HG23	1.99	0.44
1:B:219:ILE:HD12	1:B:219:ILE:C	2.43	0.44
1:C:151:LEU:HD23	1:C:229:MET:HE3	1.99	0.44
1:A:6:VAL:O	1:A:10:ILE:HG12	2.17	0.44
1:D:58:LYS:HB2	1:D:92:SER:HB3	1.99	0.44
1:E:170:LYS:O	1:E:171:ASN:HB2	2.16	0.44
1:C:36:LYS:O	1:C:55:ILE:HG22	2.18	0.44
1:D:163:CYS:SG	1:D:277:ILE:HD12	2.58	0.44
1:E:201:MET:HE3	1:E:203:ILE:HD11	2.00	0.44
1:C:208:ILE:HD13	1:F:297:PHE:CD2	2.53	0.44
1:D:101:GLN:O	1:D:104:GLN:HB3	2.18	0.44
1:A:124:VAL:O	1:A:128:ILE:HG13	2.18	0.44
1:D:154:GLY:HA2	1:D:231:ASN:ND2	2.33	0.44
1:F:49:GLN:HE22	1:F:83:ASN:ND2	2.01	0.44
1:A:212:GLU:HG3	1:A:220:THR:HG21	1.99	0.43
1:F:271:ARG:HD2	1:F:271:ARG:C	2.42	0.43
1:A:5:LEU:HD23	1:A:29:VAL:HG12	2.00	0.43
1:F:113:THR:HG21	1:F:120:LEU:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:PHE:O	1:B:66:GLN:HG2	2.18	0.43
1:C:90:THR:CG2	1:C:91:LYS:H	2.31	0.43
1:E:213:ARG:HD2	1:E:213:ARG:HA	1.77	0.43
1:E:271:ARG:HD2	1:E:276:ILE:HD11	2.01	0.43
1:A:248:THR:HG23	1:A:294:ALA:HB1	2.00	0.43
1:A:286:LEU:HD12	1:A:286:LEU:HA	1.87	0.43
1:C:286:LEU:HD12	1:C:286:LEU:HA	1.91	0.43
1:C:36:LYS:HD3	1:C:50:ILE:HG23	2.01	0.43
1:E:141:VAL:HG22	1:E:177:ASP:OD1	2.19	0.43
1:F:40:PHE:HD2	1:F:45:LEU:HD12	1.83	0.43
1:F:138:ILE:N	1:F:138:ILE:HD13	2.33	0.43
1:F:152:LEU:HD11	1:F:277:ILE:HD13	2.00	0.43
1:A:61:PHE:HB2	1:A:93:PHE:CD1	2.51	0.43
1:A:17:VAL:HA	1:A:112:LYS:O	2.18	0.43
1:A:42:MET:HB3	1:A:75:ILE:HG23	2.01	0.43
1:A:253:GLN:HE21	1:A:253:GLN:HB2	1.66	0.43
1:E:183:TYR:HD1	1:E:185:LEU:HD13	1.82	0.43
1:F:210:ASN:C	1:F:210:ASN:ND2	2.74	0.43
1:A:272:LYS:HD2	1:A:275:GLU:CD	2.44	0.43
1:C:55:ILE:HD11	1:C:90:THR:OG1	2.18	0.43
1:C:216:GLY:O	1:C:219:ILE:HG22	2.18	0.43
1:A:91:LYS:HD2	1:A:114:ASP:HA	2.00	0.43
1:A:243:PHE:CZ	1:A:266:PRO:HB3	2.54	0.43
1:A:250:LEU:HD23	1:A:250:LEU:N	2.22	0.43
1:B:267:ILE:CG2	1:E:270:GLY:HA2	2.48	0.43
1:E:91:LYS:HD3	1:E:116:PHE:CZ	2.54	0.43
1:F:240:THR:HG23	1:F:241:VAL:HG23	2.00	0.43
1:E:250:LEU:H	1:E:250:LEU:CD1	2.29	0.43
1:A:165:LEU:HA	1:A:165:LEU:HD12	1.77	0.42
1:B:164:ALA:O	1:B:168:ILE:HD13	2.19	0.42
1:D:250:LEU:HD22	1:D:287:LYS:HD3	2.01	0.42
1:E:56:LEU:HB3	1:E:93:PHE:CE1	2.54	0.42
1:A:13:PHE:O	1:A:17:VAL:HG22	2.20	0.42
1:C:8:GLU:O	1:C:12:GLN:HB2	2.20	0.42
1:A:304:ILE:HD11	1:F:195:GLU:HG2	2.01	0.42
1:A:306:LYS:HZ3	1:A:309:LYS:HE3	1.84	0.42
1:A:53:VAL:HG13	1:A:86:ALA:HB3	2.01	0.42
1:B:250:LEU:HD13	1:B:287:LYS:HD2	2.01	0.42
1:E:300:ASN:O	1:E:304:ILE:HG13	2.19	0.42
1:F:253:GLN:HE21	1:F:253:GLN:HB2	1.64	0.42
1:A:155:ARG:HG2	1:A:155:ARG:NH1	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:304:ILE:O	1:C:308:LYS:HG2	2.20	0.42
1:D:142:LEU:HB3	1:D:175:VAL:HB	2.02	0.42
1:D:210:ASN:ND2	1:D:213:ARG:H	2.17	0.42
1:E:38:VAL:HB	1:E:41:GLU:HG2	2.02	0.42
1:F:17:VAL:HG13	1:F:111:LEU:HB3	2.01	0.42
1:F:25:ASN:C	1:F:27:SER:N	2.76	0.42
1:F:145:VAL:HG13	1:F:172:HIS:CE1	2.54	0.42
1:A:168:ILE:HD12	1:A:172:HIS:O	2.20	0.42
1:C:196:VAL:HG12	1:F:304:ILE:HD12	2.02	0.42
1:D:183:TYR:HD2	1:D:185:LEU:CD1	2.30	0.42
1:C:7:LYS:HG3	1:C:8:GLU:N	2.34	0.42
1:C:219:ILE:HA	1:C:219:ILE:HD12	1.77	0.42
1:E:123:THR:O	1:F:186:GLY:HA2	2.20	0.42
1:B:204:ARG:HG3	2:B:361:HOH:O	2.19	0.42
1:C:210:ASN:C	1:C:210:ASN:ND2	2.78	0.42
1:D:58:LYS:HB2	1:D:92:SER:CB	2.49	0.42
1:A:292:ASN:HD22	1:A:292:ASN:C	2.28	0.41
1:E:35:LEU:HD12	1:E:35:LEU:HA	1.90	0.41
1:F:268:SER:HA	1:F:269:PRO:HD3	1.95	0.41
1:B:149:GLY:HA3	1:B:225:GLU:O	2.21	0.41
1:E:144:GLU:HB3	1:E:173:LEU:HB2	2.02	0.41
1:C:84:PRO:HA	1:C:85:PRO:HD3	1.90	0.41
1:D:103:ASN:C	1:D:105:THR:H	2.28	0.41
1:E:76:LEU:HD12	1:E:98:VAL:HG12	2.02	0.41
1:A:35:LEU:CD1	1:A:120:LEU:HD23	2.48	0.41
1:B:57:GLY:HA2	1:B:90:THR:CG2	2.51	0.41
1:D:49:GLN:HG3	1:D:49:GLN:O	2.20	0.41
1:F:103:ASN:ND2	1:F:108:VAL:H	2.03	0.41
1:D:89:LEU:HD13	1:D:93:PHE:CD2	2.55	0.41
1:D:97:THR:O	1:D:101:GLN:HG3	2.21	0.41
1:D:217:LEU:HD12	1:D:217:LEU:N	2.36	0.41
1:E:34:GLY:N	1:E:129:ASN:ND2	2.68	0.41
1:A:173:LEU:HD12	1:A:219:ILE:HA	2.03	0.41
1:A:210:ASN:C	1:A:210:ASN:HD22	2.29	0.41
1:B:306:LYS:C	1:B:308:LYS:H	2.28	0.41
1:D:32:VAL:CB	1:D:52:SER:HB3	2.49	0.41
1:E:34:GLY:CA	1:E:129:ASN:HD21	2.32	0.41
1:F:179:ALA:HB3	1:F:197:ALA:HB2	2.02	0.41
1:A:28:ASN:ND2	1:A:85:PRO:HA	2.35	0.41
1:C:37:ARG:NH2	2:C:318:HOH:O	2.54	0.41
1:C:38:VAL:HB	1:C:41:GLU:HG2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:PHE:HA	1:D:126:THR:OG1	2.21	0.41
1:D:149:GLY:HA3	1:D:225:GLU:O	2.21	0.41
1:D:163:CYS:SG	1:D:277:ILE:CD1	3.09	0.41
1:A:155:ARG:HG2	1:A:155:ARG:HH11	1.85	0.41
1:F:5:LEU:HB2	1:F:8:GLU:HG3	2.03	0.41
1:A:20:ILE:CD1	1:A:112:LYS:HE2	2.51	0.41
1:A:214:PHE:HB3	1:A:215:TYR:CD1	2.56	0.41
1:B:7:LYS:HB3	1:B:26:THR:CG2	2.49	0.41
1:B:126:THR:O	1:B:130:GLU:HG3	2.19	0.41
1:B:145:VAL:HA	1:B:172:HIS:CD2	2.56	0.41
1:C:228:LEU:HD11	1:C:263:TYR:HE1	1.86	0.41
1:D:97:THR:HA	1:D:100:LEU:HD12	2.02	0.41
1:E:25:ASN:C	1:E:27:SER:H	2.29	0.41
1:E:150:VAL:HG12	1:E:152:LEU:HD12	2.03	0.41
1:F:35:LEU:HD23	1:F:35:LEU:HA	1.87	0.41
1:F:98:VAL:O	1:F:102:VAL:HG23	2.20	0.41
1:F:232:LEU:HB3	1:F:267:ILE:HD11	2.03	0.41
1:A:300:ASN:O	1:A:304:ILE:HG13	2.21	0.41
1:B:91:LYS:HD2	1:B:114:ASP:O	2.21	0.41
1:A:198:LYS:HB3	1:A:199:LYS:H	1.46	0.40
1:B:172:HIS:HE1	1:B:278:GLU:OE2	2.04	0.40
1:F:13:PHE:C	1:F:15:ASP:H	2.29	0.40
1:A:32:VAL:O	1:A:52:SER:HB3	2.21	0.40
1:C:6:VAL:HG23	1:C:28:ASN:O	2.21	0.40
1:D:5:LEU:HB2	1:D:8:GLU:OE2	2.21	0.40
1:E:80:LEU:HD12	1:E:102:VAL:HG12	2.03	0.40
1:A:50:ILE:HB	1:A:84:PRO:HA	2.03	0.40
1:A:64:LEU:C	1:A:66:GLN:H	2.28	0.40
1:A:168:ILE:HD11	1:A:174:PHE:HB2	2.03	0.40
1:C:36:LYS:NZ	2:C:344:HOH:O	2.54	0.40
1:C:307:ARG:HH11	1:C:307:ARG:HG2	1.87	0.40
1:E:4:LEU:HD12	1:E:132:PHE:CE1	2.56	0.40
1:B:25:ASN:C	1:B:27:SER:N	2.78	0.40
1:D:7:LYS:HG3	1:D:8:GLU:N	2.36	0.40
1:D:25:ASN:ND2	1:D:107:GLN:HB3	2.37	0.40
1:A:142:LEU:CB	1:A:175:VAL:HB	2.52	0.40
1:B:268:SER:HA	1:B:269:PRO:HD3	1.93	0.40
1:E:122:PHE:HA	1:E:126:THR:OG1	2.21	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	299/312 (96%)	263 (88%)	31 (10%)	5 (2%)	7	13
1	B	300/312 (96%)	288 (96%)	12 (4%)	0	100	100
1	C	296/312 (95%)	279 (94%)	16 (5%)	1 (0%)	36	55
1	D	300/312 (96%)	265 (88%)	32 (11%)	3 (1%)	12	24
1	E	306/312 (98%)	289 (94%)	16 (5%)	1 (0%)	36	55
1	F	302/312 (97%)	285 (94%)	17 (6%)	0	100	100
All	All	1803/1872 (96%)	1669 (93%)	124 (7%)	10 (1%)	21	38

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	THR
1	A	94	THR
1	C	178	ASP
1	A	21	ASP
1	D	249	GLU
1	A	66	GLN
1	A	178	ASP
1	D	90	THR
1	D	306	LYS
1	E	178	ASP

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	265/279 (95%)	246 (93%)	19 (7%)	13	28
1	B	270/279 (97%)	257 (95%)	13 (5%)	23	46
1	C	265/279 (95%)	248 (94%)	17 (6%)	16	33
1	D	266/279 (95%)	255 (96%)	11 (4%)	27	53
1	E	271/279 (97%)	248 (92%)	23 (8%)	10	22
1	F	267/279 (96%)	254 (95%)	13 (5%)	22	45
All	All	1604/1674 (96%)	1508 (94%)	96 (6%)	17	36

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	19	LEU
1	A	48	SER
1	A	60	GLU
1	A	77	HIS
1	A	97	THR
1	A	122	PHE
1	A	137	GLN
1	A	165	LEU
1	A	168	ILE
1	A	173	LEU
1	A	187	ASN
1	A	218	GLN
1	A	234	SER
1	A	250	LEU
1	A	253	GLN
1	A	256	LEU
1	A	292	ASN
1	A	306	LYS
1	B	5	LEU
1	B	26	THR
1	B	79	LEU
1	B	114	ASP
1	B	165	LEU
1	B	166	ASP
1	B	168	ILE
1	B	173	LEU
1	B	218	GLN
1	B	235	LEU
1	B	255	LEU

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Mol	Chain	Res	Type
1	B	286	LEU
1	B	292	ASN
1	C	4	LEU
1	C	8	GLU
1	C	12	GLN
1	C	41	GLU
1	C	55	ILE
1	C	74	GLN
1	C	126	THR
1	C	137	GLN
1	C	152	LEU
1	C	166	ASP
1	C	189	LEU
1	C	219	ILE
1	C	234	SER
1	C	248	THR
1	C	249	GLU
1	C	256	LEU
1	C	286	LEU
1	D	21	ASP
1	D	24	THR
1	D	40	PHE
1	D	49	GLN
1	D	158	ILE
1	D	166	ASP
1	D	210	ASN
1	D	222	GLN
1	D	248	THR
1	D	253	GLN
1	D	271	ARG
1	E	4	LEU
1	E	8	GLU
1	E	9	LEU
1	E	35	LEU
1	E	38	VAL
1	E	41	GLU
1	E	56	LEU
1	E	74	GLN
1	E	79	LEU
1	E	131	GLN
1	E	145	VAL
1	E	155	ARG

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Mol	Chain	Res	Type
1	E	185	LEU
1	E	203	ILE
1	E	206	LEU
1	E	210	ASN
1	E	218	GLN
1	E	223	ARG
1	E	246	LEU
1	E	250	LEU
1	E	256	LEU
1	E	275	GLU
1	E	311	GLU
1	F	37	ARG
1	F	77	HIS
1	F	79	LEU
1	F	105	THR
1	F	135	VAL
1	F	138	ILE
1	F	152	LEU
1	F	210	ASN
1	F	249	GLU
1	F	253	GLN
1	F	255	LEU
1	F	264	GLU
1	F	266	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (77) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	49	GLN
1	A	66	GLN
1	A	83	ASN
1	A	104	GLN
1	A	131	GLN
1	A	137	GLN
1	A	171	ASN
1	A	172	HIS
1	A	187	ASN
1	A	210	ASN
1	A	218	GLN
1	A	222	GLN
1	A	253	GLN

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Mol	Chain	Res	Type
1	A	292	ASN
1	B	49	GLN
1	B	72	GLN
1	B	73	GLN
1	B	101	GLN
1	B	104	GLN
1	B	131	GLN
1	B	137	GLN
1	B	169	ASN
1	B	171	ASN
1	B	172	HIS
1	B	218	GLN
1	B	231	ASN
1	B	292	ASN
1	C	28	ASN
1	C	49	GLN
1	C	66	GLN
1	C	72	GLN
1	C	78	ASN
1	C	131	GLN
1	C	172	HIS
1	C	210	ASN
1	C	231	ASN
1	C	292	ASN
1	D	25	ASN
1	D	28	ASN
1	D	49	GLN
1	D	107	GLN
1	D	169	ASN
1	D	210	ASN
1	D	231	ASN
1	D	253	GLN
1	E	28	ASN
1	E	72	GLN
1	E	73	GLN
1	E	74	GLN
1	E	107	GLN
1	E	129	ASN
1	E	131	GLN
1	E	137	GLN
1	E	169	ASN
1	E	171	ASN

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Mol	Chain	Res	Type
1	E	210	ASN
1	E	218	GLN
1	E	231	ASN
1	E	292	ASN
1	F	12	GLN
1	F	14	GLN
1	F	28	ASN
1	F	49	GLN
1	F	73	GLN
1	F	77	HIS
1	F	83	ASN
1	F	103	ASN
1	F	104	GLN
1	F	107	GLN
1	F	139	HIS
1	F	169	ASN
1	F	194	GLN
1	F	210	ASN
1	F	222	GLN
1	F	253	GLN
1	F	292	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	F	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	F	233:LEU	C	234:SER	N	1.19

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	303/312 (97%)	1.05	78 (25%) 1 1	25, 65, 131, 145	0
1	B	304/312 (97%)	-0.03	7 (2%) 61 57	24, 44, 66, 100	0
1	C	300/312 (96%)	0.39	14 (4%) 36 32	26, 50, 90, 125	0
1	D	304/312 (97%)	0.92	71 (23%) 2 1	26, 53, 100, 113	0
1	E	310/312 (99%)	-0.08	8 (2%) 57 52	25, 41, 68, 108	0
1	F	306/312 (98%)	0.04	22 (7%) 21 19	24, 41, 90, 127	0
All	All	1827/1872 (97%)	0.38	200 (10%) 10 9	24, 46, 106, 145	0

All (200) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	19	LEU	6.5
1	F	307	ARG	6.3
1	A	110	ILE	5.5
1	D	70	VAL	5.4
1	D	107	GLN	5.4
1	D	108	VAL	5.3
1	A	115	PHE	5.1
1	F	250	LEU	5.1
1	E	238	GLN	4.8
1	D	250	LEU	4.7
1	A	17	VAL	4.5
1	D	63	PHE	4.5
1	F	239	THR	4.4
1	F	158	ILE	4.4
1	D	69	LEU	4.4
1	A	20	ILE	4.3
1	A	13	PHE	4.3
1	A	122	PHE	4.3
1	D	105	THR	4.3

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Mol	Chain	Res	Type	RSRZ
1	D	102	VAL	4.3
1	D	75	ILE	4.2
1	A	26	THR	4.2
1	A	6	VAL	4.2
1	A	113	THR	4.1
1	A	10	ILE	4.1
1	A	9	LEU	4.1
1	D	26	THR	3.9
1	D	79	LEU	3.9
1	A	87	ILE	3.8
1	A	111	LEU	3.8
1	A	94	THR	3.8
1	A	92	SER	3.8
1	C	311	GLU	3.8
1	D	109	PRO	3.8
1	D	68	THR	3.7
1	F	241	VAL	3.7
1	A	97	THR	3.6
1	A	88	ILE	3.6
1	F	306	LYS	3.6
1	A	305	LEU	3.6
1	A	24	THR	3.5
1	F	235	LEU	3.5
1	B	308	LYS	3.5
1	A	309	LYS	3.4
1	A	121	SER	3.4
1	A	108	VAL	3.4
1	A	69	LEU	3.4
1	D	106	TYR	3.3
1	C	69	LEU	3.3
1	A	93	PHE	3.3
1	D	46	PHE	3.3
1	A	106	TYR	3.3
1	D	19	LEU	3.2
1	A	4	LEU	3.2
1	A	250	LEU	3.2
1	A	102	VAL	3.2
1	D	39	VAL	3.2
1	D	307	ARG	3.2
1	A	118	THR	3.2
1	D	158	ILE	3.2
1	A	74	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
1	F	247	GLY	3.2
1	D	240	THR	3.2
1	D	236	GLU	3.1
1	A	99	LEU	3.1
1	F	236	GLU	3.1
1	A	109	PRO	3.1
1	A	242	THR	3.1
1	F	237	LYS	3.0
1	E	235	LEU	3.0
1	D	38	VAL	3.0
1	F	303	ALA	3.0
1	D	45	LEU	3.0
1	E	240	THR	3.0
1	C	46	PHE	3.0
1	A	86	ALA	3.0
1	A	76	LEU	3.0
1	C	250	LEU	3.0
1	D	80	LEU	2.9
1	D	40	PHE	2.9
1	C	40	PHE	2.9
1	F	243	PHE	2.9
1	A	70	VAL	2.9
1	D	67	LYS	2.9
1	D	22	GLY	2.9
1	C	75	ILE	2.9
1	A	98	VAL	2.9
1	A	5	LEU	2.9
1	D	1	MET	2.9
1	C	70	VAL	2.8
1	D	86	ALA	2.8
1	D	245	ARG	2.8
1	D	77	HIS	2.8
1	A	244	GLU	2.8
1	D	115	PHE	2.8
1	A	28	ASN	2.8
1	F	240	THR	2.8
1	A	23	HIS	2.8
1	D	99	LEU	2.8
1	A	25	ASN	2.7
1	A	117	SER	2.7
1	D	96	PRO	2.7
1	A	29	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	71	GLU	2.7
1	A	116	PHE	2.7
1	A	235	LEU	2.7
1	A	21	ASP	2.7
1	D	21	ASP	2.7
1	F	242	THR	2.7
1	D	27	SER	2.7
1	A	100	LEU	2.6
1	D	17	VAL	2.6
1	D	154	GLY	2.6
1	D	18	ASN	2.6
1	D	76	LEU	2.6
1	C	2	LYS	2.6
1	C	62	GLY	2.6
1	A	120	LEU	2.6
1	A	247	GLY	2.6
1	D	93	PHE	2.6
1	A	84	PRO	2.5
1	D	28	ASN	2.5
1	D	83	ASN	2.5
1	D	48	SER	2.5
1	D	241	VAL	2.5
1	F	157	GLY	2.5
1	F	304	ILE	2.5
1	A	104	GLN	2.5
1	D	78	ASN	2.5
1	A	75	ILE	2.5
1	D	43	LEU	2.5
1	D	64	LEU	2.5
1	D	98	VAL	2.5
1	A	303	ALA	2.5
1	D	42	MET	2.5
1	F	156	SER	2.5
1	D	20	ILE	2.4
1	D	82	LEU	2.4
1	C	43	LEU	2.4
1	C	305	LEU	2.4
1	A	304	ILE	2.4
1	D	111	LEU	2.4
1	E	311	GLU	2.4
1	F	244	GLU	2.4
1	F	248	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	157	GLY	2.4
1	D	85	PRO	2.3
1	D	100	LEU	2.3
1	A	78	ASN	2.3
1	D	242	THR	2.3
1	C	76	LEU	2.3
1	C	115	PHE	2.3
1	A	16	CYS	2.3
1	B	307	ARG	2.3
1	A	44	GLY	2.3
1	A	89	LEU	2.3
1	A	246	LEU	2.3
1	D	120	LEU	2.3
1	A	132	PHE	2.3
1	F	245	ARG	2.3
1	D	9	LEU	2.2
1	F	163	CYS	2.2
1	D	23	HIS	2.2
1	B	247	GLY	2.2
1	D	92	SER	2.2
1	D	110	ILE	2.2
1	A	85	PRO	2.2
1	A	90	THR	2.2
1	A	101	GLN	2.2
1	B	305	LEU	2.2
1	D	156	SER	2.2
1	E	236	GLU	2.2
1	A	54	ALA	2.2
1	A	68	THR	2.2
1	C	68	THR	2.2
1	A	79	LEU	2.2
1	D	305	LEU	2.2
1	D	29	VAL	2.1
1	A	114	ASP	2.1
1	D	47	SER	2.1
1	A	77	HIS	2.1
1	A	80	LEU	2.1
1	B	246	LEU	2.1
1	A	30	ILE	2.1
1	A	308	LYS	2.1
1	F	2	LYS	2.1
1	D	49	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	105	THR	2.1
1	A	31	ARG	2.1
1	E	309	LYS	2.1
1	A	55	ILE	2.1
1	D	87	ILE	2.1
1	A	302	LYS	2.1
1	D	288	HIS	2.1
1	D	89	LEU	2.0
1	B	236	GLU	2.0
1	E	270	GLY	2.0
1	D	114	ASP	2.0
1	E	250	LEU	2.0
1	A	126	THR	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.