



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2021 – 08:26 PM JST

PDB ID : 5AWF
Title : Crystal structure of SufB-SufC-SufD complex from Escherichia coli
Authors : Hirabayashi, K.; Wada, K.
Deposited on : 2015-07-03
Resolution : 2.96 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.1

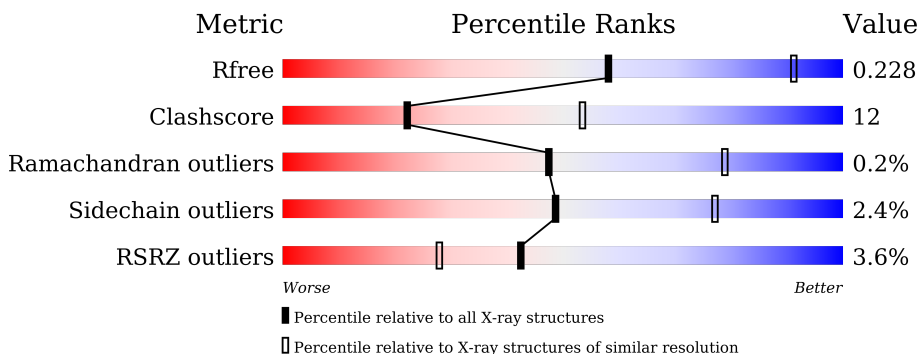
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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.



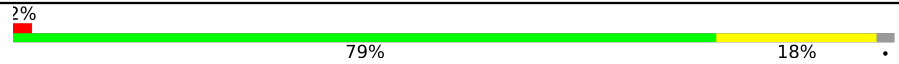

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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	495	
1	E	495	
2	B	423	
2	F	423	
3	C	248	
3	D	248	

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Mol	Chain	Length	Quality of chain
3	G	248	 <p>2% 79% 18%</p>
3	H	248	 <p>8% 59% 33% 5%</p>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 19914 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 FeS cluster assembly protein SufB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	385	Total 3004	C 1892	N 517	O 577	S 18	0	0	0
1	E	381	Total 2972	C 1872	N 511	O 571	S 18	0	0	0

- Molecule 2 is a protein called FeS cluster assembly protein SufD.

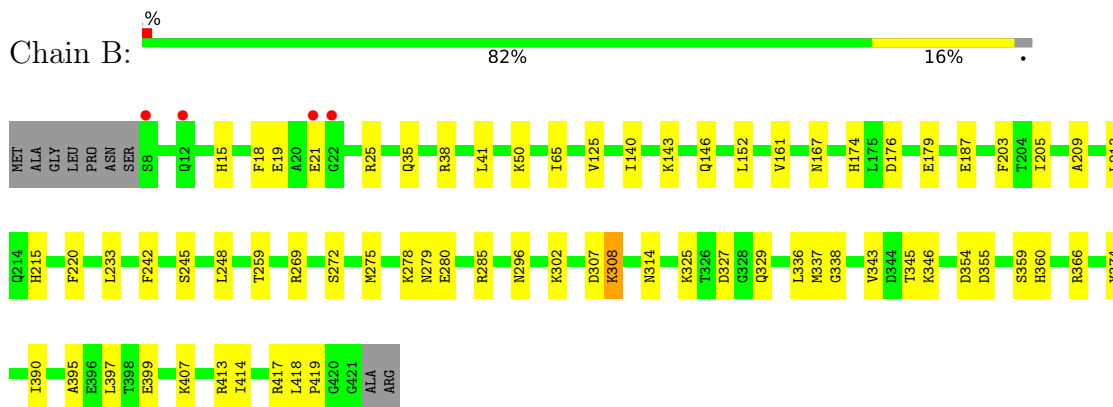
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	414	Total 3238	C 2019	N 599	O 612	S 8	0	0	0
2	F	414	Total 3238	C 2019	N 599	O 612	S 8	0	0	0

- Molecule 3 is a protein called Probable ATP-dependent transporter SufC.

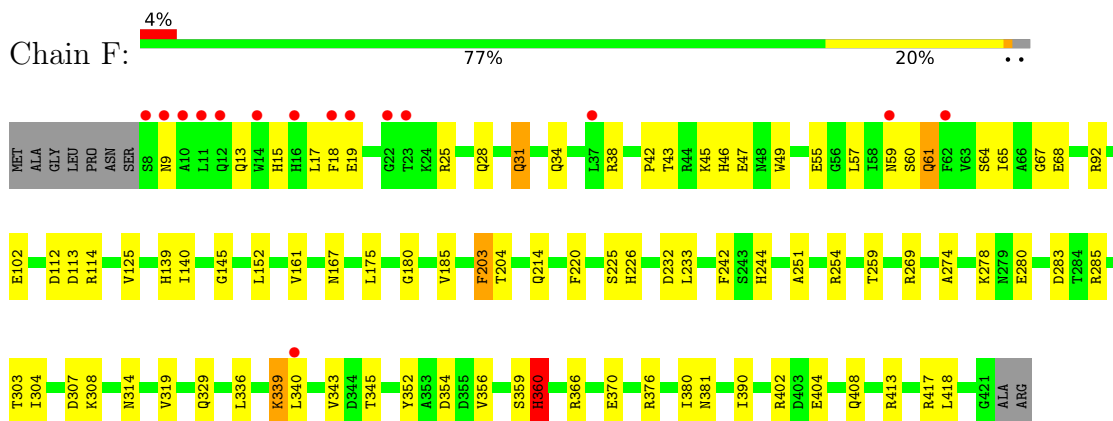
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	243	Total 1898	C 1203	N 320	O 368	S 7	0	0	0
3	D	235	Total 1829	C 1158	N 310	O 354	S 7	0	0	0
3	G	243	Total 1898	C 1203	N 320	O 368	S 7	0	0	0
3	H	236	Total 1837	C 1164	N 311	O 355	S 7	0	0	0



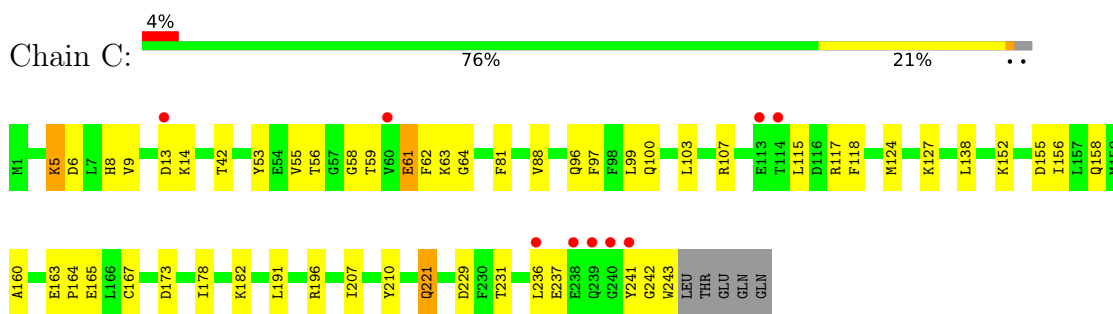
- Molecule 2: FeS cluster assembly protein SufD



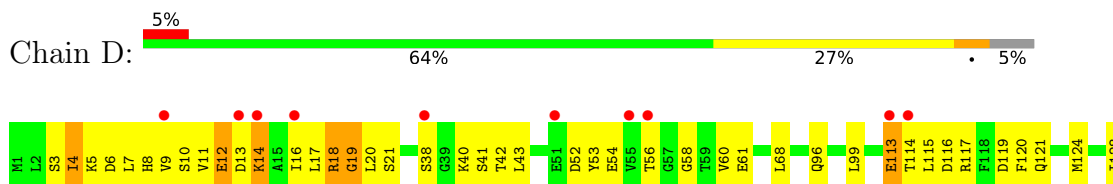
- Molecule 2: FeS cluster assembly protein SufD

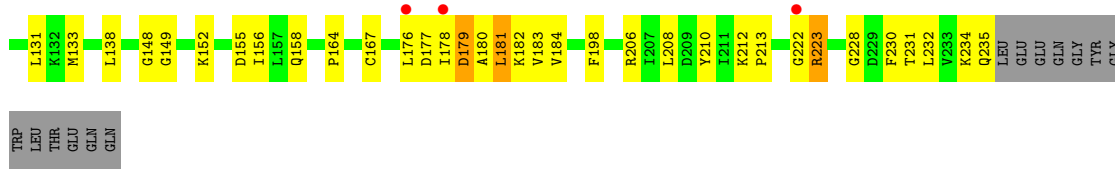


- Molecule 3: Probable ATP-dependent transporter SufC

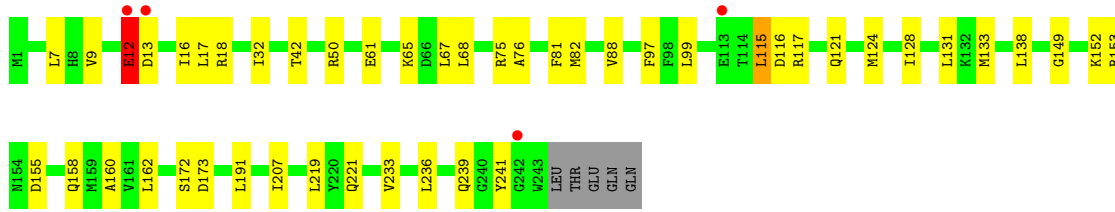
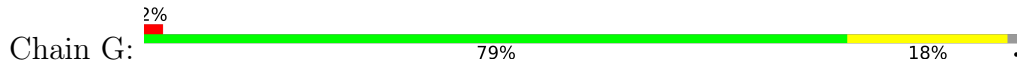


- Molecule 3: Probable ATP-dependent transporter SufC

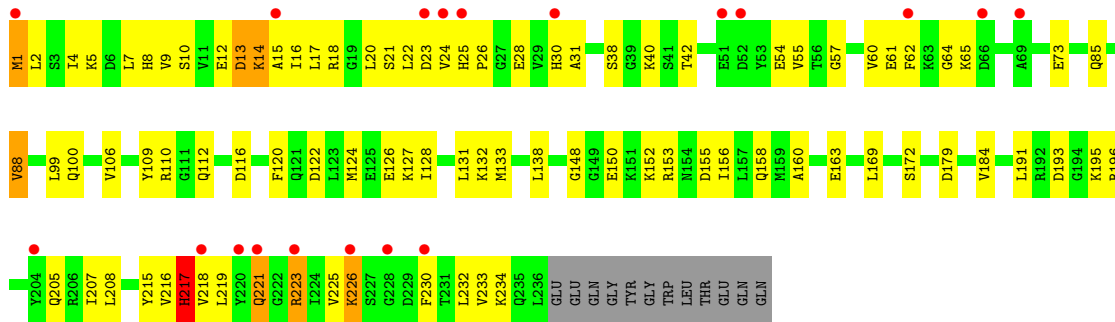




● Molecule 3: Probable ATP-dependent transporter SufC



● Molecule 3: Probable ATP-dependent transporter SufC



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.47Å 139.56Å 124.68Å 90.00° 113.10° 90.00°	Depositor
Resolution (Å)	41.11 – 2.96 43.37 – 2.96	Depositor EDS
% Data completeness (in resolution range)	98.2 (41.11-2.96) 94.7 (43.37-2.96)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.95Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.187 , 0.226 0.189 , 0.228	Depositor DCC
R_{free} test set	1995 reflections (2.58%)	wwPDB-VP
Wilson B-factor (Å ²)	63.4	Xtrriage
Anisotropy	0.541	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 71.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.015 for l,-k,h	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19914	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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 i

5.1 Standard geometry i

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	0.32	0/3068	0.58	1/4153 (0.0%)
1	E	0.31	0/3035	0.57	0/4108
2	B	0.27	0/3300	0.50	0/4473
2	F	0.30	0/3300	0.55	1/4473 (0.0%)
3	C	0.31	0/1930	0.56	0/2604
3	D	0.35	0/1858	0.65	2/2506 (0.1%)
3	G	0.30	0/1930	0.54	0/2604
3	H	0.44	0/1866	0.79	2/2517 (0.1%)
All	All	0.32	0/20287	0.59	6/27438 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	E	0	3
3	D	0	3
3	G	0	1
3	H	0	2
All	All	0	11

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	223	ARG	NE-CZ-NH2	7.01	123.80	120.30
3	D	19	GLY	N-CA-C	6.25	128.74	113.10
1	A	295	GLY	N-CA-C	6.22	128.66	113.10
3	H	14	LYS	C-N-CA	-6.17	106.28	121.70
3	H	217	HIS	N-CA-CB	5.44	120.39	110.60
2	F	360	HIS	CB-CA-C	5.32	121.03	110.40

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	294	THR	Peptide
1	A	308	GLU	Peptide
3	D	113	GLU	Peptide
3	D	12	GLU	Peptide
3	D	14	LYS	Peptide
1	E	232	THR	Peptide
1	E	290	GLY	Peptide
1	E	308	GLU	Peptide
3	G	115	LEU	Peptide
3	H	1	MET	Peptide
3	H	217	HIS	Mainchain

5.2 Too-close contacts [i](#)

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	3004	0	2916	61	0
1	E	2972	0	2887	83	0
2	B	3238	0	3184	44	0
2	F	3238	0	3184	63	0
3	C	1898	0	1896	41	0
3	D	1829	0	1840	80	0
3	G	1898	0	1896	32	0
3	H	1837	0	1851	89	0
All	All	19914	0	19654	464	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 12.

All (464) 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:E:492:HIS:CE1	1:E:493:SER:HB3	1.82	1.13
3:H:30:HIS:HD1	3:H:215:TYR:HB2	1.22	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:61:GLU:OE2	3:H:64:GLY:N	1.96	0.97
3:D:12:GLU:HG3	3:D:14:LYS:HG3	1.42	0.97
2:F:381:ASN:HA	3:H:73:GLU:HG2	1.49	0.95
3:H:30:HIS:ND1	3:H:215:TYR:HB2	1.85	0.92
3:H:217:HIS:HB3	3:H:226:LYS:O	1.70	0.90
3:D:11:VAL:O	3:D:14:LYS:N	2.04	0.89
3:D:14:LYS:HD3	3:D:16:ILE:HG12	1.56	0.88
3:H:30:HIS:CE1	3:H:217:HIS:CE1	2.63	0.86
3:D:117:ARG:HE	3:D:121:GLN:HE22	1.19	0.86
1:E:435:ALA:HB3	2:F:360:HIS:CD2	2.13	0.83
1:E:294:THR:HG23	1:E:296:GLY:H	1.43	0.83
3:D:20:LEU:N	3:D:223:ARG:HH21	1.77	0.82
3:H:124:MET:HG2	3:H:138:LEU:HD21	1.61	0.82
1:E:492:HIS:ND1	1:E:493:SER:N	2.29	0.81
3:H:16:ILE:HG21	3:H:42:THR:OG1	1.83	0.79
3:D:113:GLU:HG2	3:D:114:THR:HB	1.65	0.79
3:H:30:HIS:CE1	3:H:217:HIS:ND1	2.51	0.79
3:H:16:ILE:HD12	3:H:42:THR:HG21	1.66	0.77
1:A:320:GLU:OE1	1:A:328:LYS:NZ	2.18	0.77
2:F:67:GLY:O	2:F:92:ARG:NH2	2.18	0.76
1:A:163:GLY:HA3	1:A:212:PRO:HG3	1.66	0.76
3:H:30:HIS:NE2	3:H:217:HIS:ND1	2.33	0.76
1:E:321:THR:HG22	1:E:348:LEU:HB3	1.67	0.76
3:D:11:VAL:HG23	3:D:16:ILE:HD11	1.65	0.76
3:D:124:MET:HG2	3:D:138:LEU:HD21	1.65	0.76
3:H:223:ARG:HH11	3:H:223:ARG:HB3	1.49	0.75
2:F:42:PRO:HD3	2:F:125:VAL:HG21	1.69	0.74
3:C:6:ASP:OD1	3:C:58:GLY:HA2	1.87	0.74
3:H:223:ARG:HH11	3:H:223:ARG:CB	2.00	0.74
1:A:231:LYS:HD2	1:A:232:THR:H	1.51	0.74
3:H:223:ARG:NH1	3:H:223:ARG:HB2	2.03	0.73
3:D:10:SER:HA	3:D:16:ILE:HG13	1.70	0.73
3:H:112:GLN:NE2	3:H:163:GLU:OE2	2.20	0.72
3:H:223:ARG:CB	3:H:223:ARG:NH1	2.52	0.72
1:A:418:THR:HG22	1:A:420:PRO:HD3	1.70	0.72
1:E:289:PRO:HG3	1:E:294:THR:HA	1.70	0.72
1:A:260:ASP:HA	1:A:293:ASN:HD22	1.54	0.72
2:F:280:GLU:OE2	2:F:417:ARG:NH2	2.24	0.70
1:E:163:GLY:HA3	1:E:212:PRO:HG2	1.74	0.70
1:A:291:ASP:CG	1:A:292:ASN:H	1.95	0.69
3:G:18:ARG:NH1	3:G:221:GLN:OE1	2.25	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:214:GLN:HG2	2:F:402:ARG:HH22	1.57	0.69
2:B:35:GLN:OE1	2:B:38:ARG:NH2	2.23	0.69
3:C:5:LYS:HE2	3:C:61:GLU:HB3	1.74	0.69
3:D:18:ARG:H	3:D:223:ARG:NH1	1.89	0.69
3:H:38:SER:OG	3:H:221:GLN:N	2.25	0.69
3:D:17:LEU:HA	3:D:223:ARG:HH12	1.59	0.68
2:F:13:GLN:HE22	2:F:61:GLN:HA	1.58	0.68
3:D:8:HIS:HA	3:D:17:LEU:O	1.93	0.67
3:H:61:GLU:OE1	3:H:62:PHE:N	2.28	0.67
3:H:85:GLN:O	3:H:152:LYS:NZ	2.27	0.67
3:D:149:GLY:HA2	3:D:176:LEU:HD21	1.77	0.67
3:G:9:VAL:HG13	3:G:17:LEU:HB2	1.77	0.67
1:A:318:GLN:OE1	1:A:330:PRO:HG2	1.94	0.67
2:F:55:GLU:O	2:F:59:ASN:HB2	1.95	0.67
2:B:143:LYS:HG3	2:B:146:GLN:HG3	1.77	0.67
3:H:122:ASP:O	3:H:126:GLU:HG3	1.95	0.66
2:B:140:ILE:HD13	2:B:152:LEU:HD21	1.77	0.66
3:D:12:GLU:CG	3:D:14:LYS:HG3	2.22	0.66
2:F:46:HIS:HD1	2:F:49:TRP:HD1	1.44	0.66
2:F:204:THR:HG22	2:F:232:ASP:HB2	1.78	0.65
1:A:161:GLU:HB3	1:A:164:ILE:H	1.63	0.64
3:H:4:ILE:HG23	3:H:60:VAL:HG12	1.78	0.64
2:F:220:PHE:O	2:F:413:ARG:NH2	2.30	0.64
2:F:113:ASP:OD1	2:F:139:HIS:NE2	2.31	0.64
1:E:450:GLN:HA	1:E:450:GLN:HE21	1.62	0.63
3:H:8:HIS:HA	3:H:17:LEU:O	1.98	0.63
2:F:65:ILE:HD12	2:F:161:VAL:HG11	1.80	0.63
1:E:360:LYS:HG2	1:E:389:LEU:HB3	1.80	0.63
1:A:161:GLU:HG3	1:A:164:ILE:O	1.99	0.63
1:E:189:PRO:HD2	1:E:192:ASP:HB2	1.80	0.63
3:D:11:VAL:HB	3:D:14:LYS:HD2	1.81	0.63
3:D:11:VAL:H	3:D:16:ILE:HD11	1.64	0.63
1:E:193:ASN:OD1	1:E:194:PHE:N	2.32	0.62
1:E:294:THR:HG21	1:E:324:ALA:HA	1.80	0.62
2:F:9:ASN:OD1	2:F:13:GLN:NE2	2.32	0.62
1:A:450:GLN:OE1	1:A:451:ARG:NH1	2.31	0.62
3:C:9:VAL:HG22	3:C:55:VAL:HG22	1.82	0.62
1:E:396:ALA:O	1:E:426:ASN:ND2	2.33	0.62
3:G:115:LEU:HD12	3:G:116:ASP:HA	1.81	0.62
3:C:236:LEU:HB3	3:C:241:TYR:O	2.00	0.61
1:E:73:PRO:HB2	1:E:327:TRP:HH2	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:124:MET:HG2	3:C:138:LEU:HD21	1.83	0.61
3:H:61:GLU:OE2	3:H:65:LYS:N	2.30	0.61
3:C:237:GLU:N	3:C:237:GLU:OE1	2.34	0.60
3:H:31:ALA:HB2	3:H:208:LEU:HD21	1.83	0.60
2:F:304:ILE:HD13	2:F:390:ILE:HG23	1.83	0.60
2:B:280:GLU:OE2	2:B:417:ARG:NH2	2.34	0.60
1:A:262:TYR:HD1	1:A:294:THR:HG22	1.67	0.60
3:D:60:VAL:HB	3:D:68:LEU:HD13	1.84	0.60
2:F:404:GLU:O	2:F:408:GLN:HG2	2.02	0.60
1:E:76:LEU:HD12	1:E:77:LYS:N	2.17	0.60
3:D:156:ILE:HG13	3:D:184:VAL:HG23	1.83	0.59
3:H:1:MET:HG2	3:H:26:PRO:HD3	1.83	0.59
3:D:7:LEU:HG	3:D:9:VAL:HG23	1.84	0.59
3:D:208:LEU:HD22	3:D:213:PRO:HG3	1.83	0.59
1:E:293:ASN:O	1:E:294:THR:HG22	2.01	0.59
2:F:274:ALA:HB3	2:F:303:THR:HG22	1.84	0.59
3:C:107:ARG:NH2	3:C:163:GLU:OE2	2.35	0.59
1:E:418:THR:HG22	1:E:420:PRO:HD3	1.85	0.59
3:H:30:HIS:NE2	3:H:217:HIS:N	2.50	0.59
3:H:226:LYS:HE3	3:H:232:LEU:HD21	1.84	0.59
1:A:447:TYR:OH	3:C:155:ASP:OD1	2.20	0.59
2:B:374:TYR:OH	3:D:155:ASP:OD1	2.19	0.59
1:E:450:GLN:OE1	3:G:82:MET:HB3	2.02	0.59
3:D:99:LEU:HD21	3:D:158:GLN:HG2	1.85	0.58
1:E:294:THR:HB	1:E:353:GLN:HG2	1.85	0.58
1:E:492:HIS:HE1	1:E:493:SER:HB3	1.63	0.58
3:C:221:GLN:HE21	3:C:221:GLN:CA	2.15	0.58
1:A:296:GLY:HA3	1:A:324:ALA:HB2	1.86	0.58
1:A:262:TYR:CD1	1:A:294:THR:HG22	2.39	0.57
3:C:117:ARG:HG3	3:C:118:PHE:HD1	1.68	0.57
3:C:221:GLN:HA	3:C:221:GLN:NE2	2.19	0.57
2:F:214:GLN:HG2	2:F:402:ARG:NH2	2.19	0.57
3:D:16:ILE:HG21	3:D:42:THR:HG21	1.87	0.57
2:B:233:LEU:HB2	2:B:259:THR:HG23	1.85	0.57
3:D:11:VAL:HG22	3:D:53:TYR:HA	1.85	0.57
3:G:173:ASP:HB3	3:G:207:ILE:HD12	1.87	0.57
1:E:339:SER:HB3	1:E:368:THR:HG22	1.86	0.57
1:A:296:GLY:CA	1:A:324:ALA:HB2	2.34	0.57
3:C:221:GLN:CA	3:C:221:GLN:NE2	2.67	0.56
3:D:10:SER:O	3:D:54:GLU:HB2	2.04	0.56
3:D:38:SER:O	3:D:222:GLY:HA2	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:LYS:HD2	1:A:232:THR:N	2.19	0.56
3:H:132:LYS:HE2	3:H:153:ARG:CZ	2.34	0.56
1:E:289:PRO:HG3	1:E:294:THR:CA	2.36	0.56
1:E:293:ASN:ND2	1:E:352:HIS:HB2	2.21	0.56
3:H:150:GLU:HA	3:H:153:ARG:HH22	1.70	0.56
3:H:131:LEU:O	3:H:132:LYS:HB3	2.05	0.56
1:E:455:GLU:O	1:E:459:ILE:HG12	2.06	0.56
3:D:206:ARG:NH2	3:D:210:TYR:OH	2.37	0.56
3:D:16:ILE:HD13	3:D:42:THR:HG21	1.87	0.55
1:E:165:ILE:O	1:E:209:ILE:HA	2.05	0.55
1:E:474:LEU:HD23	1:E:479:ALA:HA	1.88	0.55
2:F:28:GLN:HA	2:F:31:GLN:CG	2.36	0.55
3:H:22:LEU:HD12	3:H:23:ASP:H	1.71	0.55
3:C:221:GLN:HE21	3:C:221:GLN:N	2.04	0.55
3:G:160:ALA:HA	3:G:191:LEU:HD21	1.88	0.55
3:H:9:VAL:HG22	3:H:55:VAL:HG22	1.87	0.55
3:C:99:LEU:HD21	3:C:158:GLN:HG2	1.88	0.55
2:F:43:THR:OG1	2:F:45:LYS:HG2	2.07	0.55
1:E:427:ASN:O	2:F:366:ARG:HB3	2.06	0.55
3:H:10:SER:O	3:H:54:GLU:HB2	2.06	0.55
3:H:128:ILE:HB	3:H:133:MET:HB2	1.88	0.55
1:A:168:SER:HB3	1:A:171:GLU:HG3	1.87	0.55
2:B:174:HIS:NE2	2:B:176:ASP:OD1	2.40	0.55
3:G:99:LEU:HD21	3:G:158:GLN:HG2	1.89	0.55
3:D:213:PRO:HD2	3:D:230:PHE:HD1	1.72	0.55
1:E:492:HIS:CG	1:E:493:SER:N	2.76	0.54
3:D:222:GLY:HA3	3:D:223:ARG:NH1	2.23	0.54
3:H:16:ILE:CD1	3:H:42:THR:HG21	2.35	0.54
3:D:213:PRO:HD2	3:D:230:PHE:CD1	2.42	0.54
1:E:363:HIS:NE2	1:E:370:SER:HB2	2.23	0.54
1:A:35:ALA:HB1	1:A:36:LYS:HA	1.90	0.54
2:F:28:GLN:HA	2:F:31:GLN:HG3	1.89	0.54
3:H:25:HIS:N	3:H:28:GLU:OE2	2.32	0.54
3:H:61:GLU:OE1	3:H:65:LYS:O	2.26	0.54
3:H:150:GLU:HA	3:H:153:ARG:NH2	2.23	0.54
1:A:433:HIS:NE2	2:B:360:HIS:HE1	2.06	0.54
1:E:356:ASP:OD2	1:E:387:ARG:NH1	2.40	0.54
3:D:179:ASP:OD2	3:D:182:LYS:HE3	2.08	0.54
3:H:1:MET:CG	3:H:25:HIS:HA	2.38	0.54
3:H:2:LEU:HB3	3:H:24:VAL:HB	1.89	0.53
2:F:339:LYS:HE3	2:F:340:LEU:HG	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:152:LYS:HD3	3:G:172:SER:O	2.08	0.53
3:H:99:LEU:HD21	3:H:158:GLN:HG2	1.89	0.53
1:A:294:THR:HB	1:A:296:GLY:CA	2.39	0.53
2:B:278:LYS:H	2:B:307:ASP:HB2	1.74	0.53
2:F:17:LEU:HD11	2:F:64:SER:CB	2.39	0.53
1:E:461:MET:HG3	3:G:97:PHE:HE2	1.74	0.53
2:F:336:LEU:HD11	2:F:343:VAL:HG12	1.90	0.53
3:D:11:VAL:HG13	3:D:52:ASP:O	2.08	0.53
3:D:18:ARG:N	3:D:223:ARG:NH1	2.57	0.52
2:F:112:ASP:OD1	2:F:114:ARG:NE	2.39	0.52
3:H:109:TYR:O	3:H:110:ARG:HB2	2.08	0.52
2:B:41:LEU:HD23	2:B:125:VAL:HG21	1.91	0.52
3:C:61:GLU:OE2	3:C:64:GLY:N	2.38	0.52
3:H:16:ILE:HG21	3:H:42:THR:CB	2.39	0.52
3:D:12:GLU:HB2	3:D:14:LYS:N	2.24	0.52
3:G:12:GLU:O	3:G:12:GLU:HG3	2.09	0.52
3:G:32:ILE:HG23	3:G:219:LEU:HD13	1.91	0.52
3:H:14:LYS:HG2	3:H:15:ALA:N	2.23	0.52
3:C:5:LYS:NZ	3:C:59:THR:HB	2.24	0.52
3:D:179:ASP:O	3:D:183:VAL:HG23	2.09	0.52
1:A:295:GLY:N	1:A:296:GLY:HA3	2.24	0.52
3:H:4:ILE:C	3:H:5:LYS:HD2	2.30	0.52
2:F:370:GLU:HG3	3:H:88:VAL:HG11	1.92	0.52
3:D:231:THR:O	3:D:232:LEU:HD12	2.09	0.51
1:A:293:ASN:ND2	1:A:294:THR:OG1	2.43	0.51
1:A:423:GLU:OE1	2:B:50:LYS:NZ	2.44	0.51
1:A:73:PRO:HB2	1:A:327:TRP:HH2	1.75	0.51
1:E:307:CYS:HB3	1:E:312:SER:HB2	1.91	0.51
2:F:46:HIS:ND1	2:F:49:TRP:HD1	2.06	0.51
1:A:301:VAL:HG11	1:A:303:LYS:HE3	1.93	0.51
1:A:375:LYS:HB3	1:A:466:PHE:CE2	2.45	0.51
1:E:308:GLU:H	1:E:309:GLY:HA3	1.76	0.51
2:B:336:LEU:HD11	2:B:343:VAL:HG12	1.92	0.51
1:E:188:VAL:O	1:E:237:ARG:NH2	2.43	0.51
1:E:174:HIS:HA	1:E:177:PRO:HG3	1.93	0.51
1:A:165:ILE:O	1:A:209:ILE:HA	2.11	0.50
1:E:161:GLU:CD	1:E:164:ILE:HG12	2.32	0.50
3:H:148:GLY:O	3:H:152:LYS:HG2	2.11	0.50
3:D:178:ILE:HD12	3:D:206:ARG:NH1	2.26	0.50
1:A:269:VAL:HB	1:A:303:LYS:HG2	1.93	0.50
3:C:160:ALA:HA	3:C:191:LEU:HD21	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:164:PRO:HG2	3:D:167:CYS:SG	2.50	0.50
3:H:223:ARG:HB2	3:H:223:ARG:CZ	2.42	0.50
1:E:456:GLU:HG2	1:E:457:ASP:N	2.26	0.50
2:F:214:GLN:NE2	2:F:244:HIS:CD2	2.79	0.50
2:F:278:LYS:H	2:F:307:ASP:HB2	1.77	0.50
3:D:11:VAL:HB	3:D:14:LYS:HZ2	1.76	0.50
3:D:116:ASP:HB3	3:D:119:ASP:OD2	2.11	0.50
1:E:258:VAL:O	1:E:289:PRO:HA	2.12	0.50
3:H:61:GLU:CD	3:H:65:LYS:H	2.14	0.50
3:D:178:ILE:HG23	3:D:206:ARG:HH12	1.77	0.50
1:A:259:ARG:O	1:A:293:ASN:ND2	2.45	0.49
2:B:161:VAL:HG13	2:B:167:ASN:HB2	1.93	0.49
3:H:127:LYS:HD2	3:H:160:ALA:HB1	1.94	0.49
3:H:132:LYS:HB3	3:H:153:ARG:HE	1.76	0.49
3:H:205:GLN:CD	3:H:205:GLN:H	2.16	0.49
1:A:439:ARG:NH1	2:B:354:ASP:OD2	2.45	0.49
3:C:117:ARG:HG3	3:C:118:PHE:CD1	2.46	0.49
3:D:11:VAL:HG12	3:D:12:GLU:H	1.76	0.49
1:E:346:VAL:HG23	1:E:470:VAL:HG11	1.94	0.49
2:B:337:MET:SD	2:B:390:ILE:HD11	2.51	0.49
1:E:238:THR:HB	1:E:269:VAL:HG22	1.94	0.49
3:H:4:ILE:HG21	3:H:20:LEU:HD22	1.95	0.49
3:H:218:VAL:HG23	3:H:226:LYS:NZ	2.28	0.49
1:E:61:LEU:O	1:E:65:ARG:HG3	2.12	0.49
1:A:42:VAL:O	1:A:46:ILE:HG13	2.13	0.49
2:B:325:LYS:HE3	2:B:355:ASP:OD1	2.12	0.49
3:D:12:GLU:HG2	3:D:14:LYS:HZ2	1.77	0.49
2:F:175:LEU:HD22	2:F:185:VAL:HG21	1.93	0.49
3:H:8:HIS:HB2	3:H:57:GLY:O	2.13	0.49
2:B:205:ILE:HG21	2:B:213:LEU:HD22	1.95	0.49
3:C:103:LEU:HD21	3:C:115:LEU:HD12	1.95	0.49
3:D:8:HIS:ND1	3:D:18:ARG:HA	2.28	0.49
1:E:450:GLN:OE1	3:G:82:MET:N	2.46	0.48
1:E:450:GLN:O	3:G:76:ALA:HB2	2.14	0.48
1:E:492:HIS:ND1	1:E:493:SER:HB3	2.25	0.48
2:F:233:LEU:HB2	2:F:259:THR:HG23	1.95	0.48
1:E:493:SER:OG	1:E:494:VAL:N	2.45	0.48
1:A:201:ALA:O	1:A:202:VAL:HB	2.13	0.48
1:A:308:GLU:N	1:A:309:GLY:HA3	2.28	0.48
2:B:395:ALA:O	2:B:399:GLU:HB2	2.12	0.48
3:D:19:GLY:C	3:D:223:ARG:HH21	2.17	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:145:GLY:HA2	2:F:180:GLY:HA3	1.96	0.48
2:B:65:ILE:HD12	2:B:161:VAL:HG11	1.95	0.48
1:E:351:GLY:O	1:E:382:SER:N	2.47	0.48
2:F:329:GLN:HA	2:F:359:SER:O	2.14	0.48
3:G:239:GLN:HB2	3:G:241:TYR:CE2	2.49	0.48
1:A:55:TRP:CD2	1:A:181:ARG:HG3	2.49	0.48
3:H:100:GLN:HG2	3:H:120:PHE:CZ	2.49	0.48
3:D:11:VAL:N	3:D:16:ILE:HD11	2.28	0.48
3:D:234:LYS:HG2	3:D:235:GLN:HG2	1.96	0.48
2:B:242:PHE:HD1	2:B:269:ARG:HB3	1.79	0.47
1:E:36:LYS:HG3	1:E:37:GLY:N	2.28	0.47
2:F:214:GLN:NE2	2:F:402:ARG:HH12	2.12	0.47
3:H:207:ILE:HG23	3:H:208:LEU:HG	1.96	0.47
1:E:294:THR:HG21	1:E:324:ALA:CA	2.44	0.47
1:E:173:ILE:O	1:E:177:PRO:HA	2.15	0.47
1:E:335:ARG:HA	1:E:364:ILE:HB	1.96	0.47
1:E:461:MET:HG3	3:G:97:PHE:CE2	2.50	0.47
3:H:12:GLU:OE1	3:H:12:GLU:HA	2.15	0.47
1:A:291:ASP:CG	1:A:292:ASN:N	2.65	0.47
3:D:117:ARG:HE	3:D:121:GLN:NE2	2.01	0.47
1:A:422:VAL:HB	2:B:345:THR:HG23	1.95	0.47
2:B:296:ASN:HA	2:B:327:ASP:O	2.14	0.47
3:D:17:LEU:HA	3:D:223:ARG:NH1	2.25	0.47
1:E:294:THR:OG1	1:E:324:ALA:N	2.47	0.47
1:E:334:LEU:HB3	1:E:339:SER:HB3	1.97	0.47
3:G:12:GLU:O	3:G:13:ASP:HB3	2.14	0.47
3:G:124:MET:HG3	3:G:138:LEU:HD11	1.96	0.47
3:H:124:MET:O	3:H:128:ILE:HG23	2.14	0.47
3:H:218:VAL:CG1	3:H:219:LEU:N	2.77	0.47
1:A:301:VAL:HG12	1:A:303:LYS:HG3	1.95	0.47
3:C:191:LEU:O	3:C:196:ARG:HD2	2.15	0.47
3:C:173:ASP:HB3	3:C:207:ILE:HD12	1.96	0.47
3:D:20:LEU:HB2	3:D:223:ARG:NH2	2.29	0.47
1:E:50:ARG:O	1:E:52:GLU:HG3	2.15	0.46
2:B:418:LEU:HB3	2:B:419:PRO:HD2	1.98	0.46
3:D:12:GLU:O	3:D:13:ASP:HB2	2.15	0.46
3:D:167:CYS:HB2	3:D:198:PHE:CD1	2.50	0.46
2:F:15:HIS:O	2:F:19:GLU:HG2	2.15	0.46
2:F:18:PHE:O	2:F:25:ARG:HD2	2.15	0.46
1:A:444:GLN:HA	3:C:88:VAL:HG11	1.98	0.46
2:B:21:GLU:H	2:B:21:GLU:HG3	1.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:16:ILE:HG21	3:H:42:THR:HG21	1.96	0.46
2:F:254:ARG:HA	2:F:283:ASP:HB3	1.98	0.46
3:G:116:ASP:OD1	3:G:117:ARG:N	2.48	0.46
2:B:278:LYS:O	2:B:279:ASN:HB2	2.16	0.46
2:B:308:LYS:HB3	2:B:308:LYS:HE3	1.50	0.46
2:F:285:ARG:HA	2:F:314:ASN:O	2.15	0.46
3:H:16:ILE:HD12	3:H:42:THR:CG2	2.40	0.46
2:B:18:PHE:O	2:B:25:ARG:NE	2.46	0.46
1:E:161:GLU:HB3	1:E:164:ILE:H	1.80	0.46
3:H:230:PHE:O	3:H:233:VAL:HG23	2.16	0.46
1:E:445:LEU:O	1:E:449:LEU:HG	2.16	0.46
3:H:110:ARG:NH2	3:H:163:GLU:O	2.49	0.46
3:H:193:ASP:OD2	3:H:196:ARG:NH1	2.49	0.46
2:B:329:GLN:HA	2:B:359:SER:O	2.17	0.45
1:A:346:VAL:HG23	1:A:470:VAL:HG11	1.98	0.45
3:D:148:GLY:O	3:D:152:LYS:HG2	2.16	0.45
3:H:153:ARG:CZ	3:H:153:ARG:HB2	2.45	0.45
1:A:475:PRO:HB2	1:A:478:PHE:HD1	1.82	0.45
3:D:12:GLU:HB2	3:D:14:LYS:H	1.81	0.45
3:D:228:GLY:HA3	3:D:232:LEU:HD11	1.97	0.45
1:A:308:GLU:H	1:A:309:GLY:HA3	1.82	0.45
1:A:445:LEU:O	1:A:449:LEU:HG	2.16	0.45
1:A:461:MET:HG3	3:C:97:PHE:HE2	1.81	0.45
2:F:68:GLU:HG2	2:F:92:ARG:HH12	1.81	0.45
2:B:15:HIS:NE2	2:B:19:GLU:OE2	2.50	0.45
2:B:399:GLU:OE2	2:B:407:LYS:NZ	2.39	0.45
1:E:308:GLU:N	1:E:309:GLY:HA3	2.32	0.45
3:H:156:ILE:HG13	3:H:184:VAL:HG13	1.98	0.45
3:C:152:LYS:O	3:C:156:ILE:HD12	2.16	0.45
3:D:14:LYS:HD3	3:D:16:ILE:CG1	2.38	0.45
3:H:216:VAL:HB	3:H:232:LEU:HD23	1.98	0.45
1:A:162:GLN:O	1:A:164:ILE:HG13	2.17	0.45
3:C:124:MET:CG	3:C:138:LEU:HD21	2.47	0.45
3:D:128:ILE:HG23	3:D:133:MET:HB2	1.99	0.45
2:F:102:GLU:OE1	2:F:102:GLU:N	2.49	0.45
2:F:278:LYS:HB2	2:F:278:LYS:HE2	1.70	0.45
3:H:191:LEU:O	3:H:196:ARG:HD2	2.17	0.45
1:A:266:ALA:HA	1:A:300:PHE:O	2.16	0.45
1:A:318:GLN:NE2	1:A:328:LYS:NZ	2.64	0.45
1:A:360:LYS:HG2	1:A:389:LEU:HB3	1.99	0.45
2:B:220:PHE:O	2:B:413:ARG:NH2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:8:HIS:O	3:D:56:THR:N	2.49	0.45
2:F:17:LEU:HD11	2:F:64:SER:HB2	1.98	0.45
3:H:13:ASP:N	3:H:13:ASP:OD1	2.49	0.45
3:D:3:SER:O	3:D:5:LYS:NZ	2.50	0.45
3:C:62:PHE:CE2	3:C:63:LYS:HE2	2.52	0.45
3:C:178:ILE:HG23	3:C:210:TYR:OH	2.17	0.45
1:E:450:GLN:HE21	1:E:450:GLN:CA	2.28	0.45
2:F:34:GLN:O	2:F:38:ARG:HG2	2.16	0.45
2:F:18:PHE:O	2:F:25:ARG:NH1	2.51	0.44
1:E:290:GLY:HA2	1:E:291:ASP:OD1	2.17	0.44
3:D:5:LYS:NZ	3:D:61:GLU:HG2	2.32	0.44
2:F:18:PHE:CZ	2:F:25:ARG:HD3	2.53	0.44
2:F:242:PHE:CD2	2:F:269:ARG:HB3	2.52	0.44
3:H:7:LEU:HD12	3:H:7:LEU:HA	1.73	0.44
1:E:417:HIS:HE1	2:F:352:TYR:CZ	2.36	0.44
3:H:30:HIS:CD2	3:H:30:HIS:C	2.90	0.44
1:A:484:LYS:HD3	1:A:484:LYS:HA	1.83	0.44
3:D:4:ILE:O	3:D:21:SER:HA	2.17	0.44
3:D:115:LEU:HD13	3:D:120:PHE:HA	2.00	0.44
2:B:220:PHE:CE2	2:B:413:ARG:HG2	2.52	0.44
1:E:435:ALA:CB	2:F:360:HIS:CD2	2.96	0.44
1:A:260:ASP:HA	1:A:293:ASN:ND2	2.29	0.44
1:E:39:ASN:OD1	1:E:42:VAL:HG23	2.18	0.44
1:E:439:ARG:NH1	2:F:354:ASP:OD2	2.51	0.44
3:G:68:LEU:HA	3:G:75:ARG:NH2	2.33	0.44
3:H:106:VAL:O	3:H:110:ARG:HD2	2.18	0.44
3:H:226:LYS:HE3	3:H:232:LEU:CD2	2.47	0.44
2:F:319:VAL:HG11	2:F:356:VAL:HG11	1.99	0.43
3:H:7:LEU:HG	3:H:9:VAL:HG23	1.99	0.43
1:A:173:ILE:HD13	1:A:180:VAL:HG11	2.00	0.43
1:E:304:ARG:HD2	1:E:333:ILE:HD11	2.00	0.43
3:C:127:LYS:HA	3:C:127:LYS:HD3	1.65	0.43
3:C:221:GLN:HE21	3:C:221:GLN:H	1.66	0.43
1:E:227:ILE:HD13	1:E:229:ALA:O	2.18	0.43
1:E:422:VAL:HB	2:F:345:THR:HG23	1.99	0.43
2:F:185:VAL:HG11	2:F:203:PHE:HE2	1.83	0.43
2:F:376:ARG:HA	2:F:380:ILE:O	2.19	0.43
2:B:187:GLU:OE2	2:B:215:HIS:NE2	2.41	0.43
3:C:138:LEU:HD23	3:C:138:LEU:HA	1.75	0.43
3:D:11:VAL:O	3:D:13:ASP:N	2.51	0.43
1:E:171:GLU:O	1:E:175:ASP:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:16:ILE:HG21	3:H:42:THR:CG2	2.48	0.43
3:D:177:ASP:O	3:D:181:LEU:HB2	2.18	0.43
3:H:132:LYS:HE2	3:H:153:ARG:NH1	2.33	0.43
3:H:169:LEU:HB3	3:H:172:SER:OG	2.18	0.43
3:C:229:ASP:OD2	3:C:231:THR:OG1	2.24	0.43
2:F:161:VAL:HG13	2:F:167:ASN:HB2	2.00	0.43
2:B:285:ARG:HA	2:B:314:ASN:O	2.18	0.43
3:D:42:THR:HG23	3:D:53:TYR:HD1	1.84	0.43
3:G:61:GLU:OE2	3:G:65:LYS:N	2.52	0.43
1:A:157:GLU:HB3	1:A:219:MET:SD	2.59	0.43
3:D:149:GLY:HA2	3:D:176:LEU:CD2	2.46	0.43
3:H:225:VAL:HG12	3:H:226:LYS:HG2	1.99	0.43
2:B:242:PHE:CD1	2:B:269:ARG:HB3	2.54	0.43
3:C:13:ASP:OD1	3:C:14:LYS:HG2	2.19	0.43
2:F:46:HIS:CD2	2:F:47:GLU:H	2.37	0.43
3:G:81:PHE:HE1	3:G:162:LEU:HD12	1.83	0.43
1:A:297:ILE:HG23	1:A:323:SER:HA	2.01	0.42
3:C:63:LYS:HD3	3:C:165:GLU:OE1	2.19	0.42
3:D:18:ARG:O	3:D:223:ARG:NE	2.52	0.42
3:D:96:GLN:HE22	3:D:121:GLN:NE2	2.17	0.42
1:E:235:PHE:CD1	1:E:266:ALA:HB3	2.54	0.42
1:E:353:GLN:O	1:E:382:SER:HB2	2.19	0.42
3:H:127:LYS:HA	3:H:127:LYS:HD3	1.82	0.42
3:D:212:LYS:HA	3:D:213:PRO:HD3	1.75	0.42
3:G:233:VAL:O	3:G:236:LEU:HB2	2.19	0.42
3:H:234:LYS:N	3:H:234:LYS:HD2	2.34	0.42
1:E:294:THR:HG21	1:E:324:ALA:CB	2.49	0.42
1:E:299:ASN:C	1:E:300:PHE:HD1	2.23	0.42
1:A:212:PRO:HG2	1:A:215:VAL:HG21	2.02	0.42
1:A:427:ASN:HB3	2:B:338:GLY:O	2.19	0.42
3:D:11:VAL:HB	3:D:14:LYS:NZ	2.33	0.42
2:B:275:MET:SD	2:B:414:ILE:HG23	2.59	0.42
2:B:245:SER:O	2:B:272:SER:HA	2.20	0.42
3:D:181:LEU:HA	3:D:184:VAL:HG12	2.01	0.42
1:E:294:THR:HG23	1:E:296:GLY:N	2.24	0.42
1:A:67:TRP:CE2	1:A:198:LEU:HD13	2.55	0.42
1:A:358:GLY:HA3	1:A:387:ARG:O	2.20	0.42
1:E:211:VAL:HA	1:E:212:PRO:HD3	1.80	0.42
1:E:450:GLN:OE1	3:G:82:MET:CB	2.66	0.42
3:G:236:LEU:HD23	3:G:236:LEU:HA	1.72	0.42
3:H:132:LYS:HB3	3:H:153:ARG:NE	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:390:ILE:HG21	2:F:418:LEU:HD21	2.01	0.41
3:H:24:VAL:HA	3:H:28:GLU:OE2	2.20	0.41
2:B:302:LYS:HB2	2:B:397:LEU:HD11	2.02	0.41
1:E:490:LEU:O	1:E:494:VAL:HG12	2.20	0.41
2:F:17:LEU:HD11	2:F:64:SER:HB3	2.02	0.41
3:G:17:LEU:HD23	3:G:17:LEU:HA	1.67	0.41
1:A:267:ALA:HB3	1:A:301:VAL:HG13	2.02	0.41
1:E:76:LEU:HD11	1:E:325:ILE:HG23	2.02	0.41
1:E:444:GLN:HG2	3:G:88:VAL:HG11	2.02	0.41
3:H:16:ILE:CG1	3:H:42:THR:HG21	2.51	0.41
3:H:179:ASP:N	3:H:179:ASP:OD1	2.53	0.41
2:B:179:GLU:HG3	2:B:209:ALA:HB3	2.03	0.41
1:E:298:LEU:HD23	1:E:324:ALA:HB3	2.02	0.41
3:H:15:ALA:HB1	3:H:18:ARG:HG3	2.03	0.41
3:D:6:ASP:O	3:D:58:GLY:HA3	2.20	0.41
1:E:446:PHE:CZ	3:G:50:ARG:HG3	2.56	0.41
2:F:242:PHE:HD2	2:F:269:ARG:HB3	1.85	0.41
3:H:4:ILE:O	3:H:21:SER:HA	2.20	0.41
1:A:427:ASN:O	2:B:366:ARG:HB3	2.20	0.41
3:C:81:PHE:HB2	3:C:164:PRO:HB3	2.03	0.41
3:C:96:GLN:O	3:C:100:GLN:HB2	2.21	0.41
1:E:446:PHE:CE2	3:G:50:ARG:HG3	2.55	0.41
3:G:149:GLY:O	3:G:153:ARG:HG3	2.20	0.41
1:A:421:TYR:CD2	2:B:346:LYS:HD2	2.56	0.41
1:E:66:ALA:O	1:E:70:MET:HG3	2.21	0.41
3:G:121:GLN:O	3:G:124:MET:HB2	2.20	0.41
3:H:8:HIS:CD2	3:H:18:ARG:HG2	2.56	0.41
2:F:57:LEU:HD13	2:F:226:HIS:CG	2.57	0.40
2:F:225:SER:O	2:F:251:ALA:N	2.44	0.40
3:H:152:LYS:O	3:H:155:ASP:HB2	2.21	0.40
1:A:39:ASN:OD1	1:A:42:VAL:HG23	2.21	0.40
3:C:5:LYS:CE	3:C:59:THR:HB	2.51	0.40
3:C:8:HIS:HB2	3:C:56:THR:O	2.21	0.40
3:C:117:ARG:NE	3:C:118:PHE:HE1	2.19	0.40
3:C:164:PRO:HG2	3:C:167:CYS:SG	2.62	0.40
3:D:18:ARG:H	3:D:223:ARG:CZ	2.34	0.40
3:D:180:ALA:O	3:D:184:VAL:HG12	2.21	0.40
3:D:181:LEU:HA	3:D:184:VAL:CG1	2.51	0.40
3:D:212:LYS:HB2	3:D:212:LYS:HE2	1.89	0.40
2:B:143:LYS:HE2	2:B:143:LYS:HB3	1.88	0.40
3:C:42:THR:HG23	3:C:53:TYR:HD1	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:242:GLY:HA2	3:C:243:TRP:HA	1.82	0.40
3:D:42:THR:HG23	3:D:53:TYR:CD1	2.56	0.40
1:E:450:GLN:HG2	1:E:451:ARG:NH1	2.37	0.40
3:G:16:ILE:HG21	3:G:42:THR:OG1	2.21	0.40
3:G:128:ILE:HG23	3:G:133:MET:HB2	2.03	0.40
3:D:40:LYS:O	3:D:43:LEU:HB3	2.22	0.40
1:E:36:LYS:CG	1:E:37:GLY:N	2.85	0.40
2:F:308:LYS:HG3	2:F:340:LEU:HD13	2.02	0.40
1:A:439:ARG:HD2	2:B:355:ASP:HA	2.03	0.40
3:D:20:LEU:HD12	3:D:222:GLY:O	2.21	0.40
2:F:140:ILE:HG21	2:F:152:LEU:HD22	2.03	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	381/495 (77%)	364 (96%)	15 (4%)	2 (0%)	29	64
1	E	377/495 (76%)	367 (97%)	10 (3%)	0	100	100
2	B	412/423 (97%)	408 (99%)	4 (1%)	0	100	100
2	F	412/423 (97%)	402 (98%)	9 (2%)	1 (0%)	47	79
3	C	241/248 (97%)	235 (98%)	6 (2%)	0	100	100
3	D	233/248 (94%)	223 (96%)	10 (4%)	0	100	100
3	G	241/248 (97%)	232 (96%)	8 (3%)	1 (0%)	34	69
3	H	234/248 (94%)	226 (97%)	8 (3%)	0	100	100
All	All	2531/2828 (90%)	2457 (97%)	70 (3%)	4 (0%)	47	79

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	202	VAL
1	A	289	PRO
2	F	60	SER
3	G	12	GLU

5.3.2 Protein sidechains [i](#)

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	323/413 (78%)	316 (98%)	7 (2%)	52 79
1	E	320/413 (78%)	308 (96%)	12 (4%)	33 66
2	B	344/350 (98%)	341 (99%)	3 (1%)	78 91
2	F	344/350 (98%)	339 (98%)	5 (2%)	65 85
3	C	207/212 (98%)	203 (98%)	4 (2%)	57 81
3	D	201/212 (95%)	195 (97%)	6 (3%)	41 72
3	G	207/212 (98%)	202 (98%)	5 (2%)	49 77
3	H	202/212 (95%)	193 (96%)	9 (4%)	27 61
All	All	2148/2374 (90%)	2097 (98%)	51 (2%)	49 77

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	LEU
1	A	184	LEU
1	A	294	THR
1	A	297	ILE
1	A	318	GLN
1	A	357	THR
1	A	493	SER
2	B	203	PHE
2	B	248	LEU
2	B	308	LYS
3	C	5	LYS
3	C	61	GLU

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Mol	Chain	Res	Type
3	C	182	LYS
3	C	221	GLN
3	D	4	ILE
3	D	18	ARG
3	D	41	SER
3	D	131	LEU
3	D	179	ASP
3	D	181	LEU
1	E	46	ILE
1	E	54	GLU
1	E	76	LEU
1	E	164	ILE
1	E	178	GLU
1	E	205	ASP
1	E	223	THR
1	E	230	GLU
1	E	414	CYS
1	E	450	GLN
1	E	492	HIS
1	E	493	SER
2	F	31	GLN
2	F	61	GLN
2	F	203	PHE
2	F	339	LYS
2	F	360	HIS
3	G	7	LEU
3	G	12	GLU
3	G	67	LEU
3	G	131	LEU
3	G	155	ASP
3	H	13	ASP
3	H	40	LYS
3	H	88	VAL
3	H	116	ASP
3	H	195	LYS
3	H	217	HIS
3	H	221	GLN
3	H	223	ARG
3	H	226	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	318	GLN
2	B	360	HIS
3	C	221	GLN
3	D	121	GLN
1	E	199	ASN
1	E	417	HIS
2	F	214	GLN
2	F	244	HIS
3	H	25	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 monosaccharides 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	385/495 (77%)	0.05	12 (3%) 49 32	29, 63, 127, 184	0
1	E	381/495 (76%)	0.13	16 (4%) 36 23	43, 81, 140, 198	0
2	B	414/423 (97%)	-0.20	4 (0%) 82 68	28, 61, 100, 154	0
2	F	414/423 (97%)	-0.00	15 (3%) 42 28	44, 78, 123, 176	0
3	C	243/248 (97%)	0.05	9 (3%) 41 27	43, 80, 140, 174	0
3	D	235/248 (94%)	0.20	13 (5%) 25 15	44, 90, 153, 194	0
3	G	243/248 (97%)	-0.04	4 (1%) 72 55	46, 80, 128, 163	0
3	H	236/248 (95%)	0.52	19 (8%) 12 6	47, 114, 176, 266	0
All	All	2551/2828 (90%)	0.06	92 (3%) 42 28	28, 78, 143, 266	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	293	ASN	6.5
3	D	222	GLY	5.6
1	A	295	GLY	5.5
1	E	229	ALA	5.2
1	E	492	HIS	4.8
3	C	239	GLN	4.5
1	E	292	ASN	4.4
2	B	8	SER	4.0
3	D	178	ILE	3.8
2	F	10	ALA	3.8
3	C	240	GLY	3.8
3	H	1	MET	3.6
2	F	23	THR	3.6
1	E	75	TRP	3.5
2	F	22	GLY	3.5
1	A	36	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	22	GLY	3.3
3	G	13	ASP	3.3
1	E	495	GLY	3.3
2	B	21	GLU	3.3
3	H	220	TYR	3.2
1	E	294	THR	3.2
2	F	37	LEU	3.2
1	A	296	GLY	3.2
3	H	23	ASP	3.2
2	F	19	GLU	3.2
2	F	62	PHE	3.2
2	F	8	SER	3.1
3	G	242	GLY	3.1
3	C	236	LEU	3.1
2	F	14	TRP	3.1
3	C	238	GLU	3.0
1	E	37	GLY	3.0
3	D	38	SER	2.9
1	E	46	ILE	2.9
3	D	113	GLU	2.9
3	D	16	ILE	2.9
3	H	226	LYS	2.9
3	H	218	VAL	2.8
2	F	59	ASN	2.8
3	H	221	GLN	2.7
2	F	11	LEU	2.7
3	C	114	THR	2.7
3	D	51	GLU	2.7
1	A	35	ALA	2.7
3	D	56	THR	2.6
3	C	13	ASP	2.6
2	F	9	ASN	2.6
1	A	71	GLU	2.6
3	H	66	ASP	2.6
3	C	241	TYR	2.6
3	G	113	GLU	2.6
1	A	294	THR	2.5
1	E	290	GLY	2.5
2	F	12	GLN	2.5
1	A	258	VAL	2.4
3	D	13	ASP	2.4
1	E	232	THR	2.4

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Mol	Chain	Res	Type	RSRZ
3	D	114	THR	2.4
3	H	204	TYR	2.4
3	H	230	PHE	2.4
3	H	62	PHE	2.4
1	A	495	GLY	2.4
3	D	9	VAL	2.3
2	F	340	LEU	2.3
3	D	55	VAL	2.3
1	E	230	GLU	2.3
3	H	30	HIS	2.3
3	H	51	GLU	2.3
1	E	257	PRO	2.2
2	F	16	HIS	2.2
3	H	15	ALA	2.2
2	B	12	GLN	2.2
3	H	223	ARG	2.2
3	H	52	ASP	2.2
1	E	36	LYS	2.2
3	D	14	LYS	2.2
3	H	69	ALA	2.2
1	A	245	ASP	2.2
3	G	12	GLU	2.2
2	F	18	PHE	2.1
3	D	176	LEU	2.1
3	C	60	VAL	2.1
3	H	24	VAL	2.1
3	H	228	GLY	2.1
3	H	25	HIS	2.1
1	E	494	VAL	2.1
3	C	113	GLU	2.1
1	A	46	ILE	2.1
1	E	295	GLY	2.0
1	A	54	GLU	2.0
1	A	49	LYS	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 monosaccharides 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.