



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 27, 2026 – 05:41 PM EDT

PDB ID : 4JFB / pdb_00004jfb
Title : Crystal structure of OmpF in C2 with tNCS
Authors : Wiseman, B.; Kilburg, A.; Chaptal, V.; Reyes-Meija, G.C.; Sarwan, J.; Falson, P.; Jault, J.M.
Deposited on : 2013-02-28
Resolution : 3.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

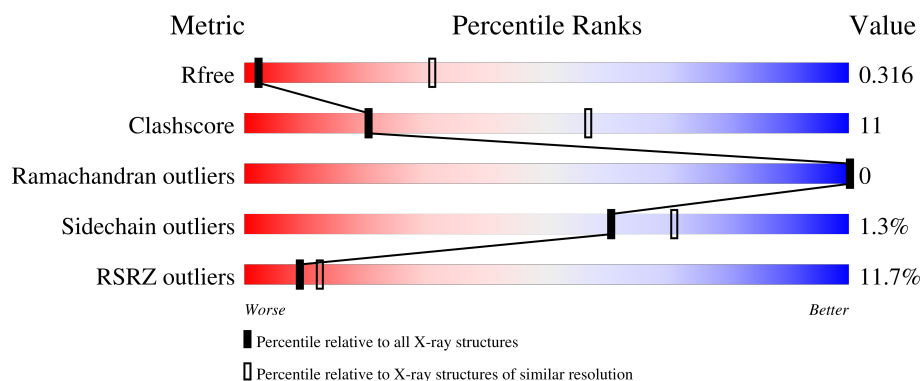
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	340	
1	B	340	
1	C	340	
1	D	340	
1	E	340	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	340	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 15765 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.

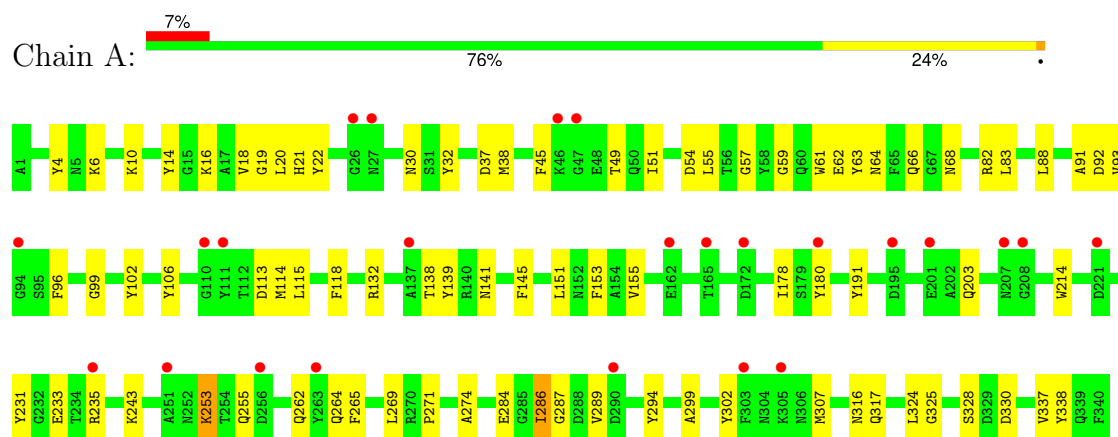
- Molecule 1 is a protein called Outer membrane protein F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2627	1654	438	532	3			
1	B	340	Total	C	N	O	S	0	0	0
			2627	1654	438	532	3			
1	C	340	Total	C	N	O	S	0	0	0
			2627	1654	438	532	3			
1	D	340	Total	C	N	O	S	0	0	0
			2627	1654	438	532	3			
1	E	340	Total	C	N	O	S	0	1	0
			2630	1656	438	533	3			
1	F	340	Total	C	N	O	S	0	0	0
			2627	1654	438	532	3			

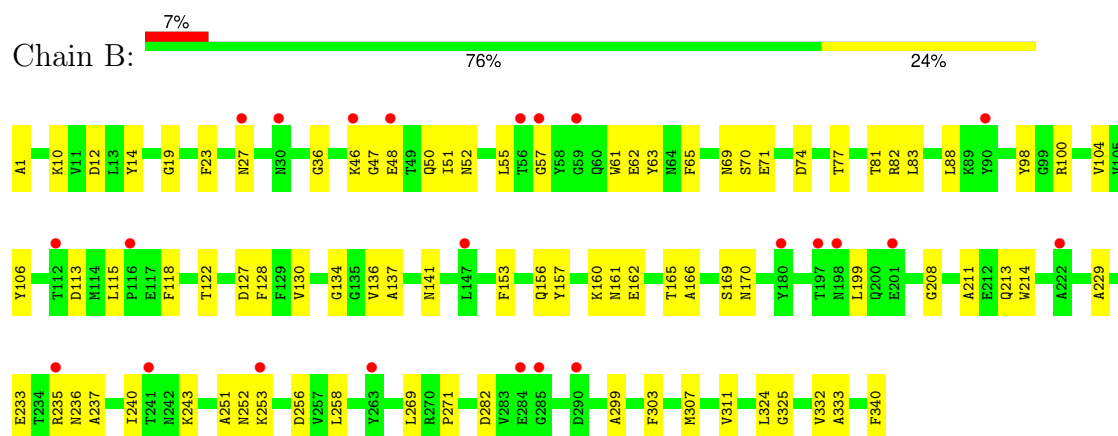
3 Residue-property plots [i](#)

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.

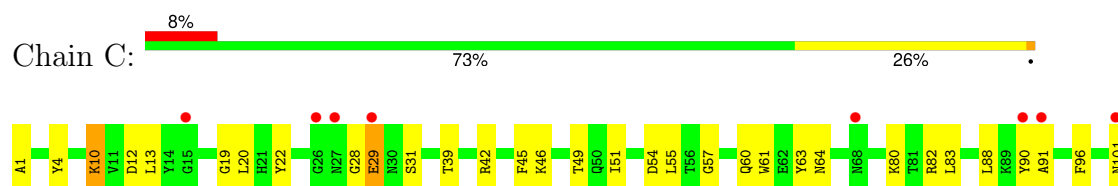
• Molecule 1: Outer membrane protein F

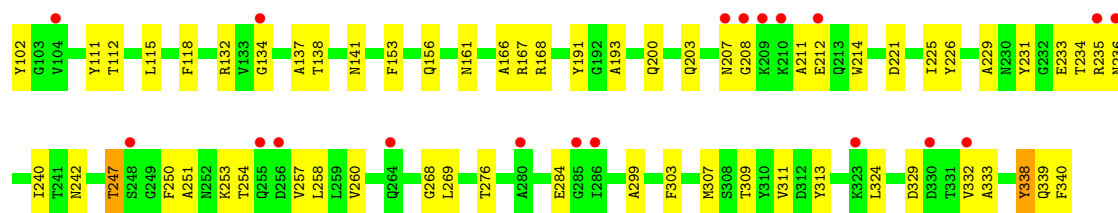


• Molecule 1: Outer membrane protein F

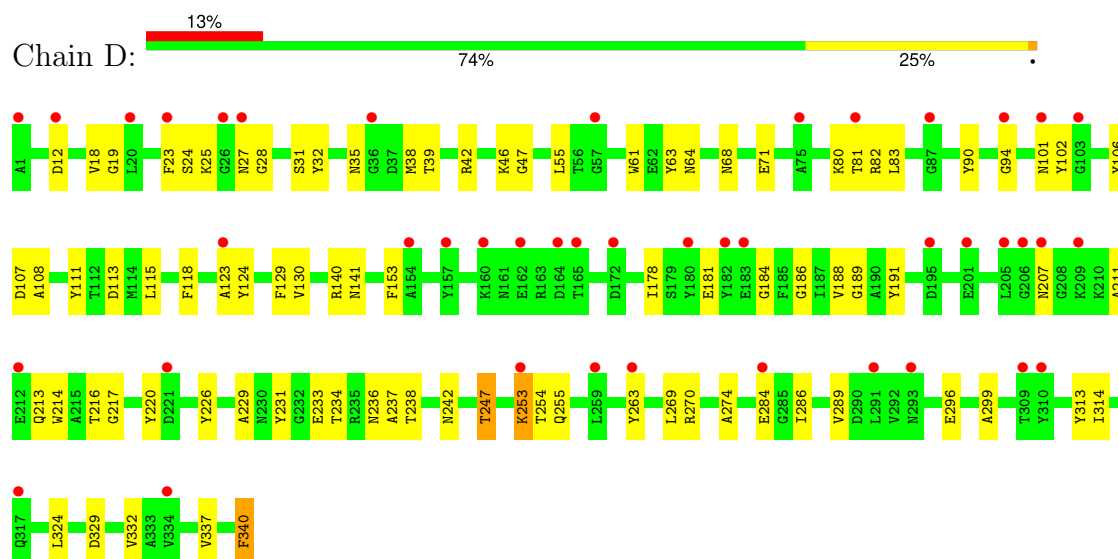


• Molecule 1: Outer membrane protein F

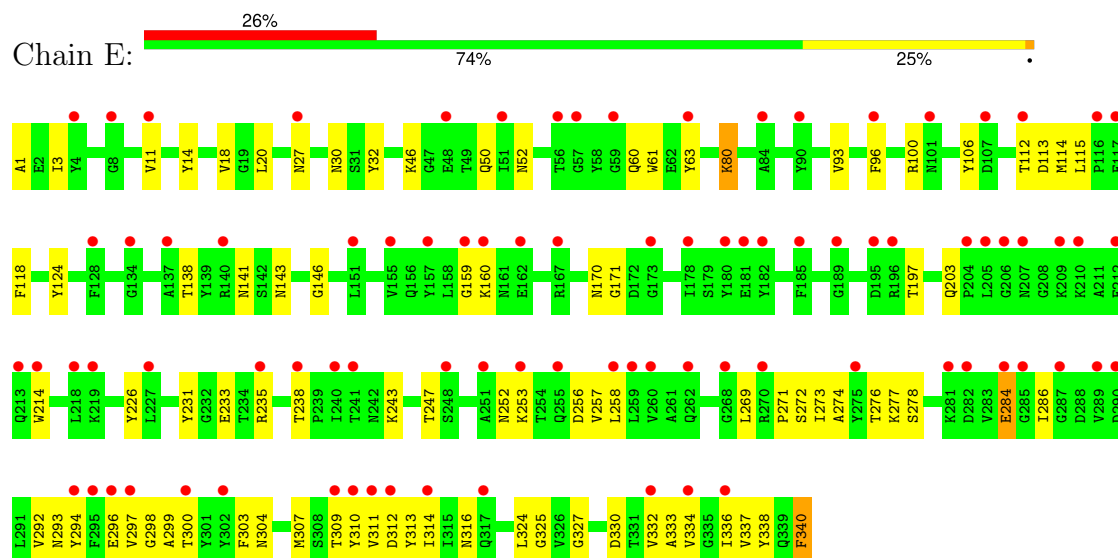




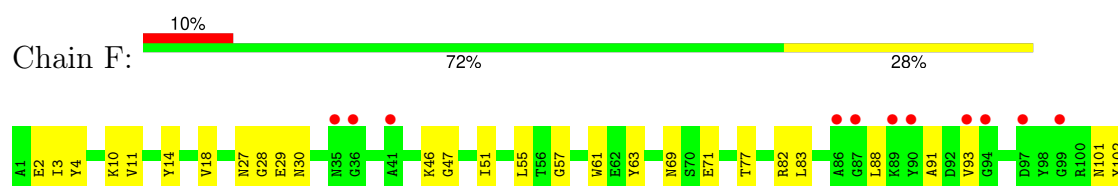
• Molecule 1: Outer membrane protein F

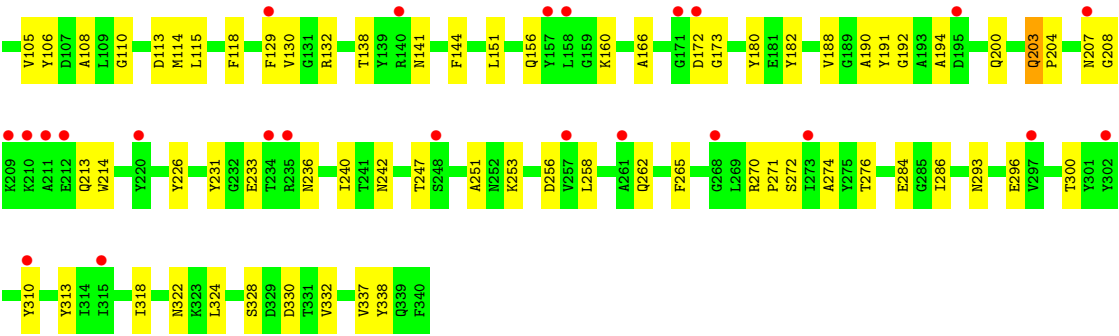


• Molecule 1: Outer membrane protein F



• Molecule 1: Outer membrane protein F





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.91Å 110.88Å 226.14Å 90.00° 104.46° 90.00°	Depositor
Resolution (Å)	68.57 – 3.80 68.57 – 3.80	Depositor EDS
% Data completeness (in resolution range)	83.3 (68.57-3.80) 83.6 (68.57-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.41 (at 3.78Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.262 , 0.314 0.263 , 0.316	Depositor DCC
R_{free} test set	1583 reflections (4.11%)	wwPDB-VP
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 129.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.77	EDS
Total number of atoms	15765	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.85 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4444e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.46	0/2683	0.88	5/3628 (0.1%)
1	B	0.44	0/2683	0.89	4/3628 (0.1%)
1	C	0.46	0/2683	0.87	7/3628 (0.2%)
1	D	0.40	0/2683	0.85	1/3628 (0.0%)
1	E	0.39	0/2690	0.81	4/3638 (0.1%)
1	F	0.42	0/2683	0.85	5/3628 (0.1%)
All	All	0.43	0/16105	0.86	26/21778 (0.1%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	29	GLU	N-CA-C	6.62	118.50	111.28
1	E	203	GLN	CA-C-N	5.93	125.14	118.97
1	E	203	GLN	C-N-CA	5.93	125.14	118.97
1	F	270	ARG	CA-C-N	5.90	125.66	119.76
1	F	270	ARG	C-N-CA	5.90	125.66	119.76
1	A	19	GLY	N-CA-C	-5.77	102.95	111.03
1	A	203	GLN	CA-C-N	5.56	124.75	118.97
1	A	203	GLN	C-N-CA	5.56	124.75	118.97
1	C	339	GLN	N-CA-C	5.40	117.39	109.24
1	C	203	GLN	CA-C-N	5.39	125.03	119.05
1	C	203	GLN	C-N-CA	5.39	125.03	119.05
1	B	36	GLY	N-CA-C	5.33	118.50	111.52
1	C	338	TYR	CA-CB-CG	-5.32	104.33	113.90
1	C	268	GLY	N-CA-C	5.29	119.24	113.58
1	F	203	GLN	CA-C-N	5.28	124.91	119.05
1	F	203	GLN	C-N-CA	5.28	124.91	119.05
1	A	16	LYS	N-CA-C	5.28	116.86	108.79
1	B	19	GLY	N-CA-C	-5.16	103.81	111.03
1	E	80	LYS	N-CA-C	5.12	116.30	108.52
1	A	287	GLY	N-CA-C	5.10	116.36	111.67
1	F	91	ALA	CB-CA-C	-5.09	110.31	117.23

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	ASN	N-CA-C	-5.09	107.20	113.41
1	C	10	LYS	N-CA-C	5.08	117.35	108.76
1	D	129	PHE	N-CA-C	5.08	119.46	113.16
1	E	27	ASN	N-CA-C	-5.06	107.24	113.41
1	B	128	PHE	CB-CA-C	-5.01	110.41	117.23

There are no chirality outliers.

There are no planarity outliers.

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	2627	0	2444	64	0
1	B	2627	0	2444	52	0
1	C	2627	0	2444	66	0
1	D	2627	0	2444	69	0
1	E	2630	0	2447	57	0
1	F	2627	0	2444	60	0
All	All	15765	0	14667	341	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 11.

All (341) 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:118:PHE:HD2	1:E:312:ASP:OD2	1.26	1.15
1:E:118:PHE:CD2	1:E:312:ASP:OD2	2.05	1.08
1:D:27:ASN:O	1:D:35:ASN:HB3	1.61	1.01
1:B:100:ARG:HG2	1:B:134:GLY:HA2	1.53	0.89
1:D:27:ASN:C	1:D:35:ASN:HB3	1.98	0.89
1:D:25:LYS:HD2	1:D:329:ASP:CG	1.99	0.88
1:E:309:THR:HG22	1:E:336:ILE:HG12	1.57	0.87
1:A:99:GLY:HA2	1:C:39:THR:HG21	1.58	0.86
1:E:160:LYS:HG3	1:E:171:GLY:HA2	1.57	0.85

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:27:ASN:O	1:D:35:ASN:CB	2.25	0.85
1:D:25:LYS:HD2	1:D:329:ASP:OD2	1.77	0.83
1:D:234:THR:HG1	1:D:254:THR:HG1	1.26	0.82
1:D:27:ASN:N	1:D:28:GLY:HA2	1.95	0.79
1:F:27:ASN:HB2	1:F:29:GLU:H	1.47	0.79
1:D:27:ASN:O	1:D:35:ASN:N	2.17	0.78
1:B:23:PHE:HB2	1:B:332:VAL:HB	1.66	0.76
1:D:27:ASN:O	1:D:35:ASN:CA	2.34	0.76
1:E:309:THR:CG2	1:E:336:ILE:HG12	2.17	0.74
1:C:29:GLU:C	1:C:31:SER:H	1.97	0.72
1:B:69:ASN:ND2	1:B:77:THR:O	2.17	0.71
1:E:93:VAL:O	1:E:141:ASN:ND2	2.24	0.70
1:A:151:LEU:HD13	1:A:180:TYR:CE2	2.27	0.70
1:A:64:ASN:HB2	1:A:82:ARG:HD2	1.72	0.69
1:B:166:ALA:HB2	1:B:199:LEU:HG	1.75	0.68
1:B:141:ASN:ND2	1:B:153:PHE:CZ	2.53	0.68
1:C:54:ASP:HB3	1:C:91:ALA:HB2	1.74	0.67
1:F:207:ASN:O	1:F:236:ASN:ND2	2.27	0.67
1:A:191:TYR:HD1	1:A:214:TRP:HB3	1.61	0.66
1:D:269:LEU:HD11	1:D:299:ALA:HB1	1.76	0.65
1:C:313:TYR:CD1	1:C:332:VAL:HG22	2.32	0.65
1:F:242:ASN:HB3	1:F:247:THR:HG22	1.78	0.65
1:E:272:SER:N	1:E:298:GLY:O	2.25	0.64
1:F:69:ASN:ND2	1:F:77:THR:O	2.21	0.64
1:A:64:ASN:ND2	1:A:66:GLN:OE1	2.30	0.64
1:D:141:ASN:HB3	1:D:153:PHE:CE1	2.34	0.63
1:C:83:LEU:HD11	1:C:102:TYR:HE1	1.63	0.63
1:A:269:LEU:HG	1:A:271:PRO:HD3	1.80	0.63
1:A:141:ASN:HB3	1:A:153:PHE:CE1	2.33	0.62
1:A:286:ILE:HG21	1:A:289:VAL:HG21	1.81	0.62
1:C:28:GLY:O	1:C:31:SER:OG	2.16	0.62
1:A:253:LYS:HE3	1:A:255:GLN:HE21	1.64	0.62
1:C:64:ASN:HB2	1:C:82:ARG:HD2	1.82	0.62
1:E:269:LEU:HD11	1:E:299:ALA:HB1	1.82	0.62
1:E:14:TYR:CE1	1:E:46:LYS:HG3	2.35	0.61
1:C:141:ASN:HB3	1:C:153:PHE:CE1	2.36	0.60
1:A:93:VAL:O	1:A:141:ASN:ND2	2.34	0.60
1:C:29:GLU:C	1:C:31:SER:N	2.59	0.60
1:F:27:ASN:HB2	1:F:29:GLU:N	2.14	0.60
1:F:27:ASN:N	1:F:28:GLY:HA2	2.16	0.60
1:A:307:MET:HG2	1:B:88:LEU:HD21	1.84	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:SER:HA	1:C:329:ASP:HB2	1.83	0.59
1:D:27:ASN:C	1:D:35:ASN:CB	2.70	0.59
1:B:1:ALA:HB1	1:C:4:TYR:CD1	2.37	0.59
1:D:24:SER:HB2	1:D:35:ASN:HB2	1.83	0.59
1:D:38:MET:HA	1:D:68:ASN:HD22	1.67	0.59
1:F:93:VAL:O	1:F:141:ASN:ND2	2.35	0.59
1:E:300:THR:HG23	1:E:310:TYR:HB3	1.85	0.59
1:E:3:ILE:HG21	1:F:3:ILE:HD12	1.85	0.59
1:D:274:ALA:HB3	1:D:296:GLU:HB3	1.85	0.58
1:D:25:LYS:CD	1:D:329:ASP:OD2	2.49	0.58
1:B:1:ALA:N	1:B:12:ASP:OD1	2.35	0.58
1:C:19:GLY:HA2	1:C:39:THR:HG22	1.84	0.57
1:F:106:TYR:O	1:F:110:GLY:N	2.35	0.57
1:A:49:THR:HG21	1:C:307:MET:HB2	1.86	0.57
1:D:12:ASP:OD1	1:D:46:LYS:HB3	2.05	0.57
1:A:51:ILE:HD12	1:A:55:LEU:HG	1.86	0.57
1:E:274:ALA:HB3	1:E:296:GLU:HB3	1.86	0.57
1:C:1:ALA:N	1:C:12:ASP:OD1	2.32	0.57
1:E:214:TRP:NE1	1:E:233:GLU:HB2	2.20	0.57
1:E:277:LYS:HA	1:E:292:VAL:O	2.04	0.57
1:A:4:TYR:CE1	1:A:6:LYS:HB3	2.39	0.57
1:B:70:SER:OG	1:B:74:ASP:OD1	2.11	0.56
1:C:101:ASN:OD1	1:C:102:TYR:N	2.37	0.56
1:D:106:TYR:OH	1:D:113:ASP:OD2	2.17	0.56
1:E:313:TYR:CD1	1:E:332:VAL:HG22	2.40	0.56
1:F:160:LYS:N	1:F:172:ASP:OD1	2.38	0.56
1:F:27:ASN:H	1:F:28:GLY:HA2	1.69	0.56
1:F:313:TYR:CD1	1:F:332:VAL:HG22	2.39	0.56
1:C:55:LEU:HD13	1:C:90:TYR:HD1	1.69	0.56
1:D:63:TYR:HD1	1:D:81:THR:HA	1.71	0.56
1:C:28:GLY:O	1:C:31:SER:CB	2.54	0.55
1:B:141:ASN:HD22	1:B:153:PHE:HZ	1.44	0.55
1:E:269:LEU:HG	1:E:271:PRO:HD3	1.88	0.55
1:E:124:TYR:HE2	1:E:238:THR:HG23	1.72	0.55
1:F:300:THR:HG23	1:F:310:TYR:HB3	1.88	0.55
1:C:212:GLU:HB2	1:C:235:ARG:HB2	1.88	0.55
1:E:316:ASN:N	1:E:330:ASP:OD1	2.38	0.55
1:A:37:ASP:OD1	1:B:98:TYR:OH	2.23	0.55
1:B:1:ALA:HB1	1:C:4:TYR:CG	2.41	0.55
1:D:80:LYS:HD3	1:F:71:GLU:HB3	1.88	0.55
1:D:24:SER:N	1:D:35:ASN:OD1	2.25	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:340:PHE:CD2	1:E:11:VAL:HG21	2.42	0.54
1:E:114:MET:HG3	1:E:226:TYR:CE2	2.43	0.54
1:A:57:GLY:HA2	1:A:88:LEU:HD23	1.89	0.54
1:A:32:TYR:HB2	1:A:316:ASN:HD21	1.71	0.54
1:E:231:TYR:HD1	1:E:257:VAL:HG22	1.73	0.54
1:A:59:GLY:HA3	1:C:338:TYR:HB3	1.90	0.54
1:B:269:LEU:HD11	1:B:299:ALA:HB1	1.89	0.54
1:D:101:ASN:OD1	1:D:102:TYR:N	2.39	0.54
1:D:42:ARG:NE	1:D:64:ASN:OD1	2.38	0.54
1:F:274:ALA:HB3	1:F:296:GLU:HB3	1.89	0.54
1:C:207:ASN:O	1:C:236:ASN:ND2	2.40	0.53
1:D:253:LYS:HE3	1:D:255:GLN:HE21	1.73	0.53
1:C:12:ASP:HB3	1:C:46:LYS:HB2	1.89	0.53
1:B:57:GLY:HA2	1:B:88:LEU:HD23	1.91	0.53
1:B:63:TYR:HD2	1:B:65:PHE:CE2	2.27	0.53
1:F:83:LEU:HD11	1:F:102:TYR:HE1	1.73	0.53
1:C:214:TRP:NE1	1:C:233:GLU:HB2	2.23	0.53
1:E:106:TYR:OH	1:E:113:ASP:OD2	2.24	0.53
1:D:207:ASN:O	1:D:236:ASN:ND2	2.41	0.53
1:A:30:ASN:ND2	1:A:328:SER:OG	2.42	0.53
1:B:214:TRP:NE1	1:B:233:GLU:HB2	2.23	0.53
1:B:160:LYS:NZ	1:B:162:GLU:OE2	2.25	0.52
1:E:159:GLY:O	1:E:170:ASN:ND2	2.43	0.52
1:B:307:MET:HB2	1:C:49:THR:HG21	1.92	0.52
1:E:61:TRP:CZ2	1:E:63:TYR:HB2	2.45	0.52
1:E:256:ASP:OD1	1:E:278:SER:OG	2.19	0.52
1:A:139:TYR:HD2	1:A:155:VAL:HG12	1.74	0.52
1:B:50:GLN:NE2	1:B:52:ASN:O	2.33	0.52
1:B:208:GLY:HA3	1:B:236:ASN:ND2	2.25	0.52
1:C:61:TRP:CZ2	1:C:63:TYR:HB2	2.45	0.52
1:D:189:GLY:HA2	1:D:216:THR:HG22	1.91	0.51
1:F:18:VAL:HG22	1:F:337:VAL:HG22	1.92	0.51
1:F:114:MET:HG3	1:F:226:TYR:CE2	2.45	0.51
1:A:115:LEU:HD21	1:A:274:ALA:HB2	1.91	0.51
1:F:138:THR:OG1	1:F:156:GLN:NE2	2.40	0.51
1:D:61:TRP:CZ2	1:D:63:TYR:HB2	2.46	0.51
1:E:303:PHE:HB2	1:E:307:MET:HB3	1.93	0.51
1:A:45:PHE:CD1	1:C:340:PHE:HE2	2.28	0.51
1:A:286:ILE:CG2	1:A:289:VAL:HG21	2.39	0.51
1:B:340:PHE:HE2	1:C:45:PHE:CD1	2.29	0.51
1:D:27:ASN:N	1:D:28:GLY:CA	2.73	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:311:VAL:HA	1:E:333:ALA:O	2.11	0.51
1:D:108:ALA:HA	1:D:188:VAL:HG11	1.93	0.50
1:E:18:VAL:HG13	1:E:337:VAL:HG22	1.93	0.50
1:D:18:VAL:HG13	1:D:337:VAL:HG22	1.93	0.50
1:A:83:LEU:HD11	1:A:102:TYR:HE1	1.77	0.50
1:C:115:LEU:HB2	1:C:118:PHE:O	2.11	0.50
1:D:211:ALA:HB1	1:D:237:ALA:HB2	1.94	0.50
1:F:101:ASN:OD1	1:F:102:TYR:N	2.40	0.50
1:A:286:ILE:CG2	1:A:289:VAL:CG2	2.89	0.50
1:A:61:TRP:HZ2	1:A:63:TYR:HD2	1.60	0.50
1:E:243:LYS:HD2	1:E:325:GLY:O	2.12	0.50
1:D:47:GLY:HA3	1:F:338:TYR:CE2	2.46	0.50
1:E:160:LYS:HE2	1:E:197:THR:HG22	1.93	0.50
1:A:30:ASN:O	1:A:328:SER:N	2.44	0.49
1:A:253:LYS:HE3	1:A:255:GLN:NE2	2.25	0.49
1:F:57:GLY:HA2	1:F:88:LEU:HD23	1.93	0.49
1:E:112:THR:HG21	1:E:258:LEU:HB2	1.94	0.49
1:E:235:ARG:HD3	1:E:253:LYS:HG2	1.94	0.49
1:E:273:ILE:HA	1:E:296:GLU:O	2.12	0.49
1:B:243:LYS:HD2	1:B:325:GLY:O	2.13	0.49
1:D:82:ARG:HA	1:F:71:GLU:OE2	2.12	0.49
1:D:107:ASP:CG	1:D:140:ARG:HH12	2.21	0.49
1:A:38:MET:HA	1:A:68:ASN:HD22	1.77	0.49
1:C:20:LEU:H	1:C:39:THR:HG22	1.78	0.49
1:A:264:GLN:HE22	1:A:302:TYR:HD2	1.59	0.49
1:F:214:TRP:NE1	1:F:233:GLU:HB2	2.28	0.49
1:A:243:LYS:HD2	1:A:325:GLY:O	2.12	0.49
1:A:338:TYR:CE2	1:B:47:GLY:HA3	2.47	0.49
1:D:231:TYR:OH	1:D:233:GLU:OE2	2.29	0.49
1:A:235:ARG:HE	1:A:253:LYS:HG3	1.77	0.48
1:B:165:THR:O	1:B:169:SER:OG	2.20	0.48
1:E:46:LYS:HG2	1:E:60:GLN:HG3	1.94	0.48
1:C:242:ASN:HA	1:C:324:LEU:HD22	1.95	0.48
1:B:252:ASN:HB3	1:B:282:ASP:HB2	1.94	0.48
1:E:340:PHE:CD2	1:F:11:VAL:HG21	2.48	0.48
1:F:115:LEU:HB2	1:F:118:PHE:O	2.14	0.48
1:C:111:TYR:O	1:C:226:TYR:OH	2.23	0.48
1:D:83:LEU:HD21	1:D:102:TYR:CD1	2.49	0.48
1:D:83:LEU:HD21	1:D:102:TYR:CE1	2.48	0.48
1:D:289:VAL:HG21	1:D:324:LEU:HG	1.96	0.48
1:D:253:LYS:HE3	1:D:255:GLN:NE2	2.29	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:VAL:HG12	1:A:20:LEU:HB2	1.95	0.48
1:A:269:LEU:HD11	1:A:299:ALA:HB1	1.96	0.48
1:F:105:VAL:HA	1:F:190:ALA:HB1	1.95	0.47
1:B:51:ILE:HD12	1:B:55:LEU:HG	1.96	0.47
1:D:18:VAL:HG22	1:D:337:VAL:HG22	1.95	0.47
1:E:115:LEU:HB2	1:E:118:PHE:O	2.15	0.47
1:C:311:VAL:HA	1:C:333:ALA:O	2.14	0.47
1:A:51:ILE:HD13	1:C:303:PHE:HB3	1.96	0.47
1:E:143:ASN:HB2	1:E:146:GLY:HA2	1.96	0.47
1:B:106:TYR:OH	1:B:113:ASP:OD2	2.18	0.47
1:D:80:LYS:HB2	1:F:71:GLU:HG2	1.96	0.47
1:C:242:ASN:HB3	1:C:247:THR:HG22	1.97	0.47
1:D:191:TYR:HD1	1:D:214:TRP:HB3	1.78	0.47
1:F:130:VAL:HG13	1:F:213:GLN:CD	2.39	0.47
1:C:234:THR:OG1	1:C:254:THR:OG1	2.17	0.47
1:F:166:ALA:HB1	1:F:200:GLN:HG2	1.97	0.47
1:A:22:TYR:CD1	1:A:38:MET:HG3	2.50	0.47
1:F:129:PHE:CD2	1:F:192:GLY:HA3	2.50	0.47
1:B:14:TYR:CE1	1:B:46:LYS:HG3	2.50	0.46
1:C:45:PHE:O	1:C:60:GLN:HG3	2.15	0.46
1:E:307:MET:SD	1:F:57:GLY:HA3	2.55	0.46
1:C:96:PHE:CZ	1:C:137:ALA:HB1	2.50	0.46
1:C:102:TYR:HA	1:C:132:ARG:HA	1.97	0.46
1:B:115:LEU:HB2	1:B:118:PHE:O	2.16	0.46
1:C:96:PHE:HA	1:C:138:THR:O	2.15	0.46
1:F:51:ILE:HD12	1:F:55:LEU:HG	1.97	0.46
1:C:82:ARG:O	1:C:132:ARG:HD2	2.15	0.46
1:C:208:GLY:HA3	1:C:236:ASN:ND2	2.30	0.46
1:B:303:PHE:HB2	1:B:307:MET:HB3	1.98	0.46
1:D:32:TYR:HE2	1:D:118:PHE:CD1	2.34	0.46
1:A:289:VAL:HG11	1:A:324:LEU:HG	1.98	0.46
1:B:61:TRP:HA	1:B:83:LEU:O	2.16	0.46
1:B:235:ARG:HD3	1:B:253:LYS:HG2	1.98	0.46
1:F:313:TYR:OH	1:F:330:ASP:OD2	2.31	0.46
1:F:318:ILE:HD12	1:F:322:ASN:ND2	2.31	0.46
1:C:338:TYR:HE2	1:C:340:PHE:HD2	1.63	0.46
1:C:240:ILE:HD13	1:C:251:ALA:HB2	1.98	0.45
1:F:108:ALA:HA	1:F:188:VAL:HG11	1.97	0.45
1:A:243:LYS:HB2	1:A:325:GLY:HA3	1.98	0.45
1:B:63:TYR:HD1	1:B:81:THR:HA	1.82	0.45
1:C:51:ILE:HD12	1:C:55:LEU:HG	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:208:GLY:HA3	1:F:236:ASN:ND2	2.32	0.45
1:B:104:VAL:N	1:B:156:GLN:OE1	2.46	0.45
1:C:338:TYR:CD2	1:C:338:TYR:C	2.92	0.45
1:D:32:TYR:CE2	1:D:314:ILE:HG13	2.52	0.45
1:E:96:PHE:HA	1:E:138:THR:O	2.16	0.45
1:F:144:PHE:N	1:F:151:LEU:O	2.49	0.45
1:B:269:LEU:HG	1:B:271:PRO:HD3	1.98	0.45
1:D:242:ASN:HB3	1:D:247:THR:HG22	1.99	0.45
1:E:294:TYR:CE1	1:E:314:ILE:HD12	2.51	0.45
1:A:22:TYR:HE2	1:A:118:PHE:CE1	2.35	0.45
1:C:83:LEU:HD11	1:C:102:TYR:CE1	2.48	0.45
1:D:31:SER:HA	1:D:329:ASP:HB2	1.98	0.45
1:F:166:ALA:HB1	1:F:200:GLN:HA	1.99	0.45
1:B:127:ASP:O	1:B:130:VAL:HG22	2.17	0.44
1:D:181:GLU:HG2	1:D:186:GLY:HA2	1.99	0.44
1:A:92:ASP:O	1:A:145:PHE:HA	2.17	0.44
1:A:294:TYR:HA	1:A:317:GLN:HG2	1.98	0.44
1:D:42:ARG:HD3	1:D:64:ASN:HA	1.98	0.44
1:F:30:ASN:O	1:F:328:SER:N	2.50	0.44
1:C:231:TYR:HD1	1:C:257:VAL:HG22	1.82	0.44
1:C:258:LEU:HG	1:C:276:THR:HG23	1.98	0.44
1:A:114:MET:C	1:A:262:GLN:HE22	2.25	0.44
1:B:324:LEU:HD23	1:B:324:LEU:HA	1.83	0.44
1:F:106:TYR:OH	1:F:113:ASP:OD2	2.35	0.44
1:F:265:PHE:HE2	1:F:271:PRO:HG3	1.81	0.44
1:A:106:TYR:OH	1:A:113:ASP:OD2	2.33	0.44
1:F:61:TRP:CZ2	1:F:63:TYR:HB2	2.52	0.44
1:F:231:TYR:HA	1:F:256:ASP:O	2.17	0.44
1:D:27:ASN:O	1:D:35:ASN:O	2.35	0.44
1:A:265:PHE:HE2	1:A:271:PRO:HG3	1.81	0.44
1:E:252:ASN:ND2	1:E:284:GLU:OE2	2.50	0.43
1:A:10:LYS:HE2	1:A:10:LYS:HB3	1.86	0.43
1:E:338:TYR:OH	1:F:47:GLY:HA3	2.18	0.43
1:E:258:LEU:HG	1:E:276:THR:HG23	2.00	0.43
1:E:276:THR:O	1:E:293:ASN:HA	2.18	0.43
1:F:262:GLN:HG2	1:F:272:SER:CB	2.49	0.43
1:A:82:ARG:O	1:A:132:ARG:HD2	2.19	0.43
1:B:130:VAL:HG13	1:B:213:GLN:CD	2.43	0.43
1:B:160:LYS:HE2	1:B:162:GLU:HG3	2.01	0.43
1:C:42:ARG:HD3	1:C:64:ASN:HA	1.99	0.43
1:D:313:TYR:HD1	1:D:332:VAL:HG22	1.83	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:309:THR:HG22	1:E:336:ILE:CG1	2.40	0.43
1:D:71:GLU:HG3	1:E:100:ARG:NH2	2.33	0.43
1:D:217:GLY:HA2	1:D:229:ALA:O	2.19	0.43
1:D:313:TYR:CD1	1:D:332:VAL:HG22	2.53	0.43
1:F:82:ARG:O	1:F:132:ARG:HD2	2.19	0.43
1:A:307:MET:SD	1:A:338:TYR:HB2	2.58	0.43
1:D:18:VAL:HG22	1:D:337:VAL:HG13	2.00	0.43
1:E:50:GLN:NE2	1:E:52:ASN:O	2.37	0.43
1:A:4:TYR:HE1	1:A:6:LYS:HB3	1.84	0.43
1:F:105:VAL:HA	1:F:190:ALA:CB	2.49	0.43
1:B:62:GLU:O	1:B:82:ARG:N	2.47	0.43
1:D:141:ASN:HB3	1:D:153:PHE:HE1	1.79	0.43
1:B:61:TRP:CZ2	1:B:63:TYR:HB2	2.53	0.42
1:B:161:ASN:HB2	1:B:170:ASN:CG	2.44	0.42
1:C:20:LEU:HD21	1:C:22:TYR:HE1	1.84	0.42
1:C:191:TYR:HD1	1:C:214:TRP:HB3	1.84	0.42
1:F:180:TYR:CE2	1:F:182:TYR:HB2	2.54	0.42
1:A:54:ASP:HB3	1:A:91:ALA:HB2	2.01	0.42
1:A:286:ILE:HG22	1:A:289:VAL:CG2	2.49	0.42
1:D:184:GLY:O	1:D:220:TYR:HA	2.19	0.42
1:D:123:ALA:HA	1:D:130:VAL:O	2.20	0.42
1:C:134:GLY:O	1:C:161:ASN:ND2	2.52	0.42
1:C:138:THR:OG1	1:C:156:GLN:NE2	2.52	0.42
1:B:71:GLU:HB2	1:C:168:ARG:HH12	1.84	0.42
1:E:324:LEU:HD23	1:E:324:LEU:HA	1.94	0.42
1:A:316:ASN:HD22	1:A:330:ASP:HA	1.84	0.42
1:D:55:LEU:HD13	1:D:90:TYR:HD1	1.85	0.42
1:E:304:ASN:ND2	1:F:51:ILE:HA	2.34	0.42
1:D:115:LEU:HB2	1:D:118:PHE:O	2.19	0.42
1:D:130:VAL:HG13	1:D:213:GLN:CD	2.44	0.42
1:E:309:THR:HG22	1:E:336:ILE:HG23	2.02	0.42
1:B:122:THR:HG21	1:B:256:ASP:HB3	2.01	0.42
1:E:311:VAL:HG22	1:E:334:VAL:HG22	2.00	0.42
1:B:311:VAL:HA	1:B:333:ALA:O	2.20	0.42
1:C:193:ALA:HA	1:C:211:ALA:O	2.20	0.42
1:D:124:TYR:HE2	1:D:238:THR:HG23	1.85	0.42
1:C:269:LEU:HD11	1:C:299:ALA:HB1	2.02	0.42
1:C:1:ALA:O	1:C:13:LEU:N	2.53	0.41
1:C:57:GLY:HA2	1:C:88:LEU:HD23	2.02	0.41
1:F:2:GLU:OE1	1:F:10:LYS:HD2	2.20	0.41
1:A:21:HIS:HD2	1:A:22:TYR:O	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:TYR:OH	1:A:62:GLU:OE1	2.36	0.41
1:A:32:TYR:HE2	1:A:118:PHE:CD1	2.38	0.41
1:D:19:GLY:HA2	1:D:39:THR:HG23	2.01	0.41
1:D:263:TYR:O	1:D:270:ARG:HA	2.20	0.41
1:E:32:TYR:HE2	1:E:118:PHE:CD1	2.38	0.41
1:F:191:TYR:HD1	1:F:214:TRP:HB3	1.84	0.41
1:F:324:LEU:HD23	1:F:324:LEU:HA	1.92	0.41
1:D:111:TYR:O	1:D:226:TYR:OH	2.36	0.41
1:F:258:LEU:HG	1:F:276:THR:HG23	2.02	0.41
1:B:137:ALA:HB3	1:B:157:TYR:HB3	2.01	0.41
1:C:4:TYR:O	1:C:10:LYS:HA	2.20	0.41
1:C:166:ALA:HB1	1:C:200:GLN:HA	2.02	0.41
1:A:153:PHE:HB2	1:A:178:ILE:HG12	2.03	0.41
1:A:231:TYR:OH	1:A:233:GLU:OE2	2.39	0.41
1:E:80:LYS:HE2	1:E:80:LYS:HB3	1.93	0.41
1:F:14:TYR:CE1	1:F:46:LYS:HG3	2.55	0.41
1:A:18:VAL:HG22	1:A:337:VAL:HG22	2.02	0.41
1:B:229:ALA:HA	1:B:258:LEU:O	2.21	0.41
1:A:18:VAL:HG13	1:A:337:VAL:HG22	2.02	0.41
1:A:96:PHE:HA	1:A:138:THR:O	2.21	0.41
1:B:136:VAL:HA	1:B:157:TYR:O	2.20	0.41
1:C:221:ASP:HA	1:C:225:ILE:O	2.21	0.41
1:D:63:TYR:CD1	1:D:81:THR:HA	2.54	0.41
1:D:94:GLY:CA	1:D:141:ASN:HD22	2.34	0.41
1:E:294:TYR:CD1	1:E:314:ILE:HG23	2.55	0.41
1:F:173:GLY:HA3	1:F:194:ALA:HB2	2.02	0.41
1:B:71:GLU:HB3	1:C:80:LYS:HD2	2.02	0.41
1:C:229:ALA:HA	1:C:258:LEU:O	2.21	0.41
1:E:1:ALA:HB1	1:F:4:TYR:CD2	2.55	0.41
1:A:21:HIS:HD2	1:A:22:TYR:N	2.19	0.40
1:B:211:ALA:HB1	1:B:237:ALA:HB2	2.03	0.40
1:F:240:ILE:HD13	1:F:251:ALA:HB2	2.01	0.40
1:A:45:PHE:CE1	1:C:340:PHE:HE2	2.39	0.40
1:C:112:THR:O	1:C:260:VAL:HG21	2.22	0.40
1:D:153:PHE:HB2	1:D:178:ILE:HG12	2.03	0.40
1:E:30:ASN:OD1	1:E:327:GLY:HA2	2.21	0.40
1:F:203:GLN:HA	1:F:204:PRO:HD3	1.90	0.40
1:C:167:ARG:HA	1:C:250:PHE:CZ	2.57	0.40
1:D:23:PHE:HB2	1:D:332:VAL:HB	2.03	0.40
1:A:32:TYR:HB2	1:A:316:ASN:ND2	2.34	0.40
1:A:32:TYR:CB	1:A:316:ASN:HD21	2.32	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:LYS:HG2	1:B:48:GLU:O	2.21	0.40
1:B:240:ILE:HD13	1:B:251:ALA:HB2	2.03	0.40
1:F:276:THR:O	1:F:293:ASN:HA	2.22	0.40
1:E:20:LEU:HD13	1:E:310:TYR:OH	2.22	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	338/340 (99%)	321 (95%)	17 (5%)	0	100	100
1	B	338/340 (99%)	319 (94%)	19 (6%)	0	100	100
1	C	338/340 (99%)	320 (95%)	18 (5%)	0	100	100
1	D	338/340 (99%)	321 (95%)	17 (5%)	0	100	100
1	E	339/340 (100%)	323 (95%)	16 (5%)	0	100	100
1	F	338/340 (99%)	322 (95%)	16 (5%)	0	100	100
All	All	2029/2040 (100%)	1926 (95%)	103 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	263/263 (100%)	260 (99%)	3 (1%)	65	73
1	B	263/263 (100%)	263 (100%)	0	100	100
1	C	263/263 (100%)	259 (98%)	4 (2%)	57	69
1	D	263/263 (100%)	258 (98%)	5 (2%)	50	66
1	E	264/263 (100%)	259 (98%)	5 (2%)	50	66
1	F	263/263 (100%)	260 (99%)	3 (1%)	65	73
All	All	1579/1578 (100%)	1559 (99%)	20 (1%)	61	71

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	253	LYS
1	A	284	GLU
1	A	286	ILE
1	C	247	THR
1	C	253	LYS
1	C	284	GLU
1	C	309	THR
1	D	247	THR
1	D	253	LYS
1	D	284	GLU
1	D	286	ILE
1	D	340	PHE
1	E	247	THR
1	E	284	GLU
1	E	286	ILE
1	E	297	VAL
1	E	340	PHE
1	F	253	LYS
1	F	284	GLU
1	F	286	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	255	GLN
1	A	316	ASN
1	B	203	GLN
1	C	76	GLN
1	C	322	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	9	ASN
1	D	68	ASN
1	D	76	GLN
1	D	161	ASN
1	D	170	ASN
1	D	264	GLN
1	E	223	ASN
1	F	9	ASN
1	F	21	HIS
1	F	161	ASN
1	F	170	ASN
1	F	317	GLN
1	F	339	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	340/340 (100%)	0.74	24 (7%)	22 19	25, 91, 154, 205	1 (0%)
1	B	340/340 (100%)	0.76	23 (6%)	23 20	18, 97, 153, 214	0
1	C	340/340 (100%)	0.76	27 (7%)	18 17	24, 83, 142, 188	1 (0%)
1	D	340/340 (100%)	1.08	43 (12%)	8 11	53, 136, 193, 246	0
1	E	340/340 (100%)	1.47	87 (25%)	1 3	59, 157, 214, 267	2 (0%)
1	F	340/340 (100%)	0.96	35 (10%)	12 14	51, 112, 175, 209	1 (0%)
All	All	2040/2040 (100%)	0.96	239 (11%)	9 12	18, 109, 190, 267	5 (0%)

All (239) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	312	ASP	6.0
1	B	285	GLY	5.6
1	F	209	LYS	5.3
1	E	282	ASP	5.2
1	D	165	THR	4.9
1	D	182	TYR	4.9
1	F	212	GLU	4.7
1	B	30	ASN	4.7
1	C	209	LYS	4.5
1	E	101	ASN	4.4
1	E	285	GLY	4.4
1	E	90	TYR	4.3
1	F	210	LYS	4.3
1	D	27	ASN	4.3
1	D	207	ASN	4.3
1	A	251	ALA	4.2
1	E	180	TYR	4.2
1	E	56[A]	THR	4.1
1	F	129	PHE	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	208	GLY	4.1
1	D	291	LEU	4.0
1	C	29	GLU	4.0
1	E	112	THR	4.0
1	E	209	LYS	3.9
1	E	162	GLU	3.8
1	D	103	GLY	3.8
1	A	207	ASN	3.8
1	D	101	ASN	3.8
1	D	157	TYR	3.7
1	C	27	ASN	3.6
1	C	248	SER	3.6
1	A	46	LYS	3.6
1	C	210	LYS	3.6
1	E	57	GLY	3.6
1	E	235	ARG	3.5
1	D	309	THR	3.5
1	C	207	ASN	3.5
1	F	248	SER	3.5
1	F	172	ASP	3.5
1	E	336	ILE	3.4
1	E	241	THR	3.4
1	D	201	GLU	3.4
1	B	235	ARG	3.3
1	E	270	ARG	3.3
1	E	287	GLY	3.3
1	A	165	THR	3.3
1	F	220	TYR	3.3
1	B	201	GLU	3.3
1	E	302	TYR	3.2
1	E	240	ILE	3.2
1	D	36	GLY	3.2
1	E	157	TYR	3.2
1	F	93	VAL	3.2
1	E	210	LYS	3.1
1	A	195	ASP	3.1
1	D	94	GLY	3.1
1	E	173	GLY	3.1
1	E	314	ILE	3.1
1	D	180	TYR	3.1
1	A	208	GLY	3.1
1	F	158	LEU	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	300	THR	3.1
1	D	162	GLU	3.1
1	D	317	GLN	3.0
1	D	20	LEU	3.0
1	D	12	ASP	3.0
1	E	107	ASP	3.0
1	C	235	ARG	3.0
1	D	183	GLU	3.0
1	E	238	THR	3.0
1	F	310	TYR	3.0
1	C	264	GLN	3.0
1	E	251	ALA	3.0
1	F	157	TYR	3.0
1	E	11	VAL	3.0
1	D	195	ASP	3.0
1	D	259	LEU	2.9
1	A	110	GLY	2.9
1	C	26	GLY	2.9
1	E	253	LYS	2.9
1	D	1	ALA	2.9
1	E	332	VAL	2.9
1	B	253	LYS	2.9
1	D	212	GLU	2.9
1	F	235	ARG	2.9
1	E	134	GLY	2.9
1	E	205	LEU	2.9
1	E	185	PHE	2.8
1	E	212	GLU	2.8
1	C	134	GLY	2.8
1	E	84	ALA	2.8
1	E	189	GLY	2.8
1	E	167	ARG	2.8
1	E	181	GLU	2.8
1	E	309	THR	2.8
1	E	196	ARG	2.8
1	A	305	LYS	2.7
1	B	180	TYR	2.7
1	E	128	PHE	2.7
1	E	259	LEU	2.7
1	E	260	VAL	2.7
1	E	137	ALA	2.7
1	D	26	GLY	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	90	TYR	2.7
1	D	23	PHE	2.7
1	C	256	ASP	2.7
1	E	116	PRO	2.6
1	E	310	TYR	2.6
1	F	89	LYS	2.6
1	B	112	THR	2.6
1	B	284	GLU	2.6
1	E	27	ASN	2.6
1	E	255	GLN	2.6
1	E	159	GLY	2.6
1	F	99	GLY	2.6
1	F	171	GLY	2.6
1	A	162	GLU	2.6
1	D	87	GLY	2.6
1	F	195	ASP	2.6
1	D	75	ALA	2.6
1	A	290	ASP	2.6
1	F	207	ASN	2.6
1	E	4	TYR	2.5
1	E	195	ASP	2.5
1	A	137	ALA	2.5
1	B	46	LYS	2.5
1	D	263	TYR	2.5
1	D	334	VAL	2.5
1	B	198	ASN	2.5
1	F	97	ASP	2.5
1	C	101	ASN	2.5
1	D	209	LYS	2.5
1	F	211	ALA	2.5
1	E	227	LEU	2.5
1	E	289	VAL	2.5
1	E	151	LEU	2.5
1	D	205	LEU	2.4
1	C	323	LYS	2.4
1	B	116	PRO	2.4
1	A	47	GLY	2.4
1	C	104	VAL	2.4
1	D	253	LYS	2.4
1	E	297	VAL	2.4
1	E	117	GLU	2.4
1	E	296	GLU	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	280	ALA	2.4
1	E	206	GLY	2.4
1	B	263	TYR	2.4
1	C	255	GLN	2.3
1	C	286	ILE	2.3
1	C	332	VAL	2.3
1	E	8	GLY	2.3
1	E	59	GLY	2.3
1	F	257	VAL	2.3
1	A	263	TYR	2.3
1	E	294	TYR	2.3
1	D	154	ALA	2.3
1	A	201	GLU	2.3
1	E	48	GLU	2.3
1	E	268	GLY	2.3
1	F	297	VAL	2.3
1	A	111	TYR	2.3
1	E	262	GLN	2.3
1	F	315	ILE	2.3
1	B	59	GLY	2.3
1	E	96	PHE	2.3
1	D	81	THR	2.3
1	F	35	ASN	2.3
1	E	214	TRP	2.3
1	E	204	PRO	2.3
1	A	235	ARG	2.3
1	B	197	THR	2.2
1	D	293	ASN	2.2
1	F	140	ARG	2.2
1	B	56	THR	2.2
1	A	221	ASP	2.2
1	B	27	ASN	2.2
1	A	180	TYR	2.2
1	E	63	TYR	2.2
1	E	51	ILE	2.2
1	B	290	ASP	2.2
1	B	147	LEU	2.2
1	A	94	GLY	2.2
1	D	57	GLY	2.2
1	F	94	GLY	2.2
1	A	256	ASP	2.2
1	E	258	LEU	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	140	ARG	2.2
1	E	213	GLN	2.2
1	D	206	GLY	2.2
1	E	178	ILE	2.2
1	C	68	ASN	2.2
1	F	86	ALA	2.1
1	B	90	TYR	2.1
1	D	123	ALA	2.1
1	A	26	GLY	2.1
1	E	284	GLU	2.1
1	E	248	SER	2.1
1	E	317	GLN	2.1
1	F	273	ILE	2.1
1	D	160	LYS	2.1
1	E	281	LYS	2.1
1	B	57	GLY	2.1
1	B	48	GLU	2.1
1	C	91	ALA	2.1
1	E	155	VAL	2.1
1	E	295	PHE	2.1
1	D	172	ASP	2.1
1	D	221	ASP	2.1
1	A	27	ASN	2.1
1	D	310	TYR	2.1
1	A	172	ASP	2.1
1	C	330	ASP	2.1
1	F	261	ALA	2.1
1	E	207	ASN	2.1
1	E	311	VAL	2.1
1	E	334	VAL	2.1
1	C	15	GLY	2.0
1	E	275	TYR	2.0
1	F	87	GLY	2.0
1	F	268	GLY	2.0
1	F	302	TYR	2.0
1	F	234	THR	2.0
1	E	160	LYS	2.0
1	E	218	LEU	2.0
1	C	236	ASN	2.0
1	C	285	GLY	2.0
1	C	90	TYR	2.0
1	C	212	GLU	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	164	ASP	2.0
1	D	284	GLU	2.0
1	E	290	ASP	2.0
1	A	303	PHE	2.0
1	B	222	ALA	2.0
1	F	41	ALA	2.0
1	E	219	LYS	2.0
1	F	36	GLY	2.0
1	E	182	TYR	2.0
1	B	241	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.