Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:18 PM GMT

PDB ID	:	4RFS
Title	:	Structure of a pantothenate energy coupling factor transporter
Authors	:	Zhang, M.; Bao, Z.; Zhao, Q.; Guo, H.; Xu, K.; Zhang, P.
Deposited on	:	2014-09-27
Resolution	:	3.23 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

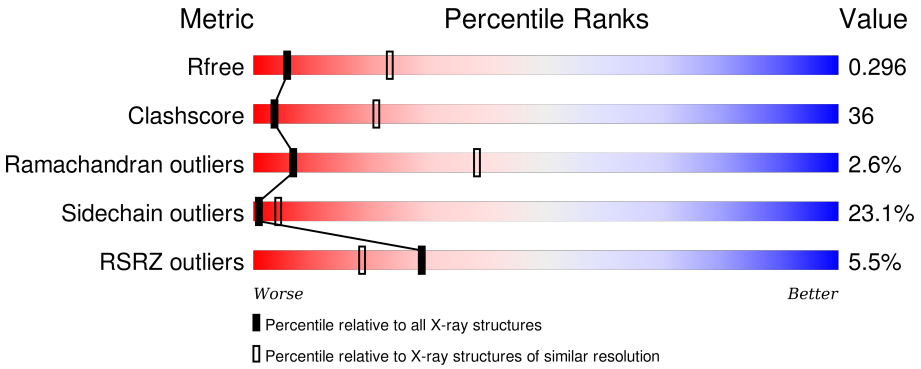
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.23 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}	91344	1092 (3.28-3.20)
Clashscore	102246	1227 (3.28-3.20)
Ramachandran outliers	100387	1204 (3.28-3.20)
Sidechain outliers	100360	1203 (3.28-3.20)
RSRZ outliers	91569	1097 (3.28-3.20)

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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	290	3% 60% 31% 7% •
2	B	279	5% 55% 32% 11% •
3	S	203	2% 36% 40% 14% • 8%
4	T	280	9% 38% 32% 13% • 16%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7601 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 Energy-coupling factor transporter ATP-binding protein EcfA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2217	1399	389	418	11			

- Molecule 2 is a protein called Energy-coupling factor transporter ATP-binding protein EcfA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	275	Total	C	N	O	S	0	0	0
			2114	1325	365	417	7			

- Molecule 3 is a protein called Substrate binding pritein S.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	S	187	Total	C	N	O	S	0	0	0
			1453	973	242	230	8			

- Molecule 4 is a protein called Energy-coupling factor transporter transmembrane protein EcfT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	T	236	Total	C	N	O	S	0	0	0
			1817	1200	295	310	12			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
T	-13	MET	-	EXPRESSION TAG	UNP Q03PY7
T	-12	GLY	-	EXPRESSION TAG	UNP Q03PY7
T	-11	SER	-	EXPRESSION TAG	UNP Q03PY7
T	-10	SER	-	EXPRESSION TAG	UNP Q03PY7
T	-9	HIS	-	EXPRESSION TAG	UNP Q03PY7
T	-8	HIS	-	EXPRESSION TAG	UNP Q03PY7
T	-7	HIS	-	EXPRESSION TAG	UNP Q03PY7

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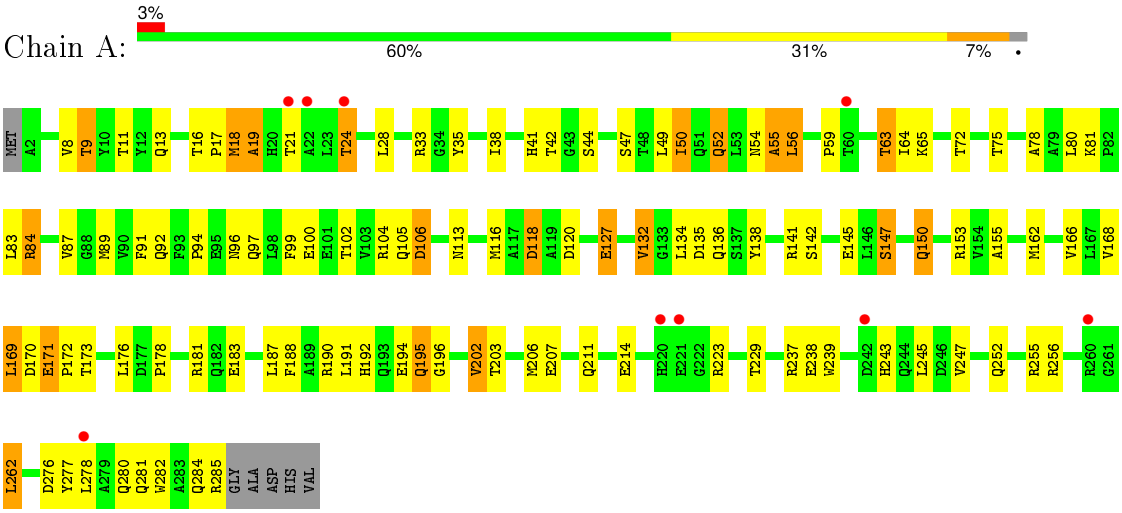
Chain	Residue	Modelled	Actual	Comment	Reference
T	-6	HIS	-	EXPRESSION TAG	UNP Q03PY7
T	-5	HIS	-	EXPRESSION TAG	UNP Q03PY7
T	-4	HIS	-	EXPRESSION TAG	UNP Q03PY7
T	-3	SER	-	EXPRESSION TAG	UNP Q03PY7
T	-2	GLN	-	EXPRESSION TAG	UNP Q03PY7
T	-1	ASP	-	EXPRESSION TAG	UNP Q03PY7
T	0	PRO	-	EXPRESSION TAG	UNP Q03PY7

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

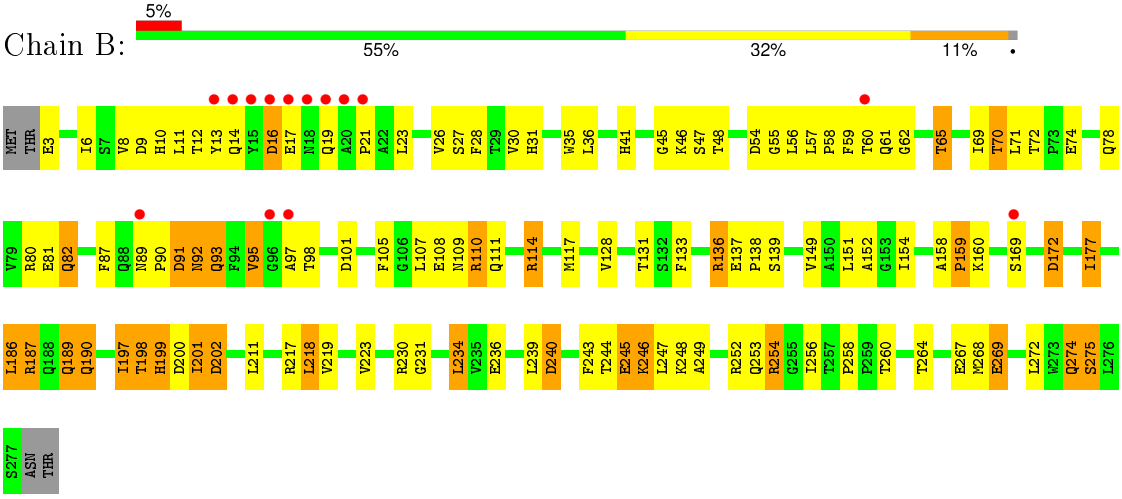
- Molecule 1: Energy-coupling factor transporter ATP-binding protein EcfA2

Chain A:



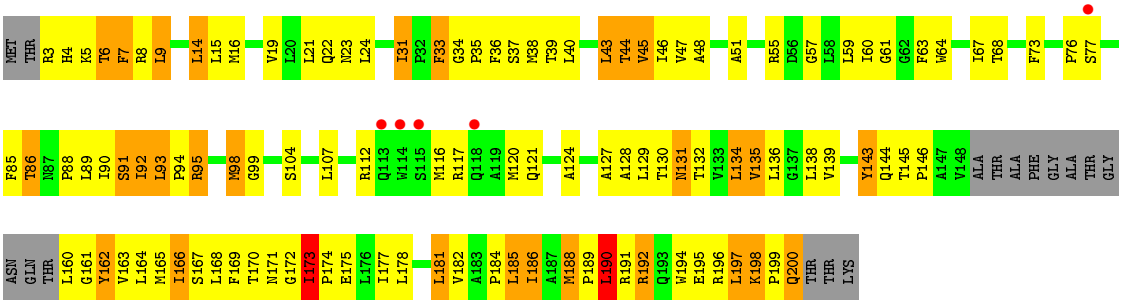
- Molecule 2: Energy-coupling factor transporter ATP-binding protein EcfA1

Chain B:



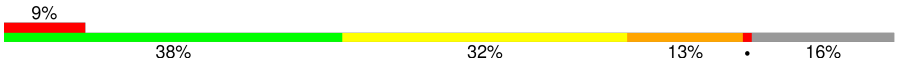
- Molecule 3: Substrate binding pritein S

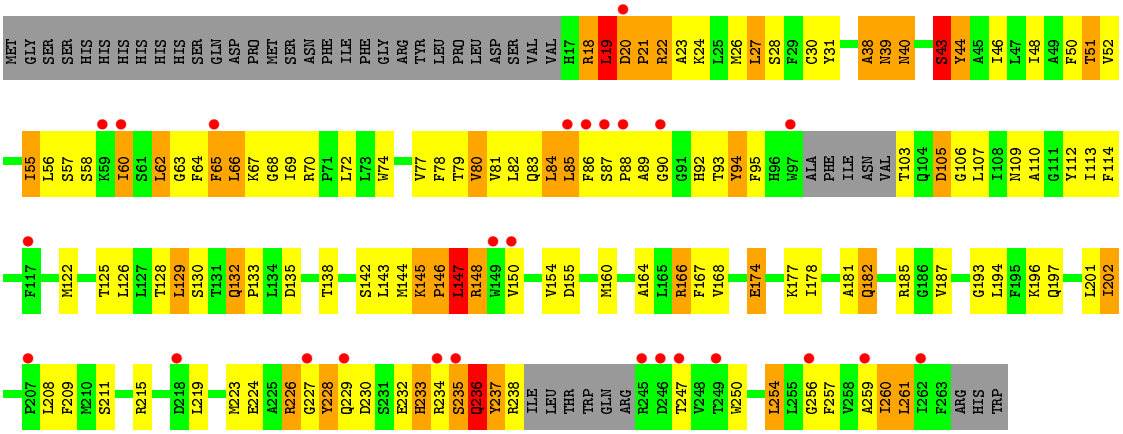
Chain S:



- Molecule 4: Energy-coupling factor transporter transmembrane protein EcfT

Chain T:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	77.73Å 145.24Å 157.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.50 – 3.23 43.97 – 3.23	Depositor EDS
% Data completeness (in resolution range)	97.1 (37.50-3.23) 90.6 (43.97-3.23)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.10 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.226 , 0.289 0.239 , 0.296	Depositor DCC
R_{free} test set	1306 reflections (4.94%)	DCC
Wilson B-factor (Å^2)	105.9	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 28321 reflections	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	7601	wwPDB-VP
Average B, all atoms (Å^2)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.58	1/2262 (0.0%)	0.73	2/3071 (0.1%)
2	B	0.48	0/2148	0.65	0/2922
3	S	0.42	0/1492	0.66	0/2038
4	T	0.42	0/1853	0.63	4/2513 (0.2%)
All	All	0.49	1/7755 (0.0%)	0.67	6/10544 (0.1%)

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
4	T	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	106	ASP	C-N	6.61	1.49	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	T	94	TYR	O-C-N	-7.79	110.23	122.70
1	A	105	GLN	O-C-N	-6.81	111.80	122.70
4	T	94	TYR	CA-C-N	5.28	128.82	117.20
4	T	94	TYR	C-N-CA	5.27	134.87	121.70
1	A	176	LEU	CA-CB-CG	5.02	126.84	115.30
4	T	261	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	T	226	ARG	Sidechain

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	2217	0	2194	83	0
2	B	2114	0	2104	98	0
3	S	1453	0	1543	208	0
4	T	1817	0	1875	210	0
All	All	7601	0	7716	551	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 36.

All (551) 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)
4:T:256:GLY:O	4:T:260:ILE:CD1	1.65	1.43
3:S:47:VAL:HA	3:S:190:LEU:CD2	1.56	1.36
4:T:63:GLY:HA2	4:T:66:LEU:CD1	1.58	1.34
3:S:143:TYR:CA	3:S:164:LEU:HG	1.58	1.32
4:T:88:PRO:HG3	4:T:103:THR:CB	1.60	1.32
4:T:65:PHE:CE2	4:T:69:ILE:HD11	1.69	1.25
3:S:160:LEU:HB3	3:S:163:VAL:CG1	1.66	1.25
4:T:63:GLY:CA	4:T:66:LEU:HD11	1.69	1.21
3:S:195:GLU:O	3:S:199:PRO:HD3	1.41	1.21
3:S:143:TYR:HA	3:S:164:LEU:CG	1.70	1.20
4:T:92:HIS:HB3	4:T:94:TYR:CD2	1.76	1.19
4:T:256:GLY:O	4:T:260:ILE:HD12	1.39	1.19
4:T:63:GLY:O	4:T:66:LEU:HD12	1.37	1.18
3:S:191:ARG:O	3:S:195:GLU:HG3	1.39	1.18
3:S:46:ILE:CG2	3:S:186:ILE:HD12	1.72	1.18
3:S:9:LEU:HD23	3:S:9:LEU:O	1.39	1.17
4:T:92:HIS:HB3	4:T:94:TYR:HD2	1.04	1.15
3:S:46:ILE:HG21	3:S:186:ILE:HD12	1.19	1.15
2:B:111:GLN:NE2	4:T:233:HIS:HD2	1.41	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:160:LEU:HB3	3:S:163:VAL:HG11	1.25	1.14
3:S:185:LEU:CD2	4:T:202:ILE:HG12	1.77	1.14
3:S:51:ALA:HB2	3:S:190:LEU:CD1	1.75	1.14
4:T:256:GLY:O	4:T:260:ILE:HD11	1.45	1.12
4:T:18:ARG:HB3	4:T:130:SER:HB2	1.19	1.12
4:T:46:ILE:HG21	4:T:260:ILE:CG1	1.79	1.11
4:T:65:PHE:HE2	4:T:69:ILE:CD1	1.63	1.11
4:T:19:LEU:HD13	4:T:19:LEU:H	1.10	1.11
3:S:51:ALA:CB	3:S:190:LEU:HD11	1.80	1.10
4:T:84:LEU:O	4:T:88:PRO:HD2	1.53	1.08
3:S:47:VAL:HA	3:S:190:LEU:HD21	1.30	1.07
4:T:46:ILE:HG21	4:T:260:ILE:HG12	1.26	1.07
4:T:46:ILE:HG21	4:T:260:ILE:CD1	1.86	1.06
3:S:161:GLY:O	3:S:165:MET:HG2	1.52	1.06
3:S:46:ILE:HB	3:S:186:ILE:HD11	1.10	1.06
3:S:143:TYR:HB2	3:S:164:LEU:HB3	1.38	1.06
3:S:46:ILE:HB	3:S:186:ILE:CD1	1.87	1.05
3:S:185:LEU:O	3:S:185:LEU:HD13	1.56	1.03
4:T:85:LEU:HD23	4:T:85:LEU:O	1.58	1.00
4:T:63:GLY:CA	4:T:66:LEU:CD1	2.30	0.99
3:S:162:TYR:HE1	4:T:90:GLY:O	1.42	0.99
3:S:143:TYR:HB2	3:S:164:LEU:CB	1.92	0.99
4:T:65:PHE:CE2	4:T:69:ILE:CD1	2.39	0.99
4:T:65:PHE:HE2	4:T:69:ILE:HD11	0.83	0.98
3:S:51:ALA:HB2	3:S:190:LEU:HD11	0.98	0.97
2:B:111:GLN:NE2	4:T:233:HIS:CD2	2.34	0.96
4:T:63:GLY:C	4:T:66:LEU:HD12	1.84	0.96
4:T:63:GLY:HA2	4:T:66:LEU:HD11	0.96	0.96
3:S:46:ILE:CB	3:S:186:ILE:HD11	1.97	0.95
4:T:62:LEU:O	4:T:66:LEU:HD11	1.66	0.95
4:T:19:LEU:N	4:T:19:LEU:CD1	2.30	0.95
3:S:182:VAL:HA	3:S:186:ILE:HB	1.46	0.95
4:T:62:LEU:O	4:T:66:LEU:HD21	1.66	0.94
4:T:19:LEU:CD1	4:T:19:LEU:H	1.80	0.94
3:S:132:THR:O	3:S:135:VAL:HG23	1.66	0.94
4:T:46:ILE:CG2	4:T:260:ILE:CD1	2.45	0.94
3:S:3:ARG:O	3:S:6:THR:HG23	1.67	0.94
3:S:185:LEU:HD21	4:T:202:ILE:HG12	1.51	0.93
4:T:60:ILE:HD12	4:T:60:ILE:O	1.67	0.93
3:S:46:ILE:CG2	3:S:186:ILE:CD1	2.46	0.93
3:S:47:VAL:HA	3:S:190:LEU:HD22	1.49	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:47:VAL:HG22	3:S:190:LEU:HD22	1.49	0.93
3:S:9:LEU:HD23	3:S:9:LEU:C	1.90	0.91
4:T:19:LEU:N	4:T:19:LEU:HD13	1.84	0.91
3:S:182:VAL:CA	3:S:186:ILE:HB	2.00	0.90
3:S:47:VAL:CA	3:S:190:LEU:CD2	2.49	0.90
3:S:134:LEU:HD23	3:S:134:LEU:N	1.85	0.90
3:S:46:ILE:CB	3:S:186:ILE:CD1	2.49	0.90
3:S:165:MET:HA	3:S:168:LEU:HD12	1.52	0.90
3:S:160:LEU:O	3:S:164:LEU:CD2	2.20	0.89
3:S:162:TYR:CE1	4:T:90:GLY:O	2.25	0.89
3:S:134:LEU:HD23	3:S:134:LEU:H	1.37	0.88
4:T:92:HIS:CG	4:T:94:TYR:HE2	1.92	0.88
4:T:105:ASP:O	4:T:109:ASN:HB2	1.74	0.88
3:S:191:ARG:O	3:S:195:GLU:CG	2.23	0.87
4:T:82:LEU:HG	4:T:86:PHE:CE2	2.11	0.86
3:S:196:ARG:O	3:S:199:PRO:HD2	1.75	0.86
1:A:194:GLU:O	1:A:195:GLN:HB3	1.76	0.86
3:S:160:LEU:O	3:S:163:VAL:HG13	1.77	0.85
4:T:92:HIS:CG	4:T:94:TYR:CE2	2.65	0.84
2:B:111:GLN:HE22	4:T:233:HIS:HD2	1.25	0.84
2:B:97:ALA:O	2:B:138:PRO:HD3	1.77	0.84
1:A:195:GLN:HG3	1:A:195:GLN:O	1.78	0.84
4:T:18:ARG:CB	4:T:130:SER:HB2	2.06	0.83
3:S:166:ILE:HG12	4:T:86:PHE:CE1	2.12	0.83
3:S:139:VAL:HG12	3:S:168:LEU:HD21	1.60	0.83
3:S:185:LEU:CD1	3:S:185:LEU:C	2.47	0.83
3:S:143:TYR:CD1	3:S:143:TYR:C	2.51	0.83
3:S:117:ARG:HE	3:S:191:ARG:HH12	1.27	0.82
3:S:196:ARG:O	3:S:199:PRO:HG2	1.78	0.82
4:T:85:LEU:C	4:T:85:LEU:CD2	2.47	0.82
3:S:143:TYR:CB	3:S:164:LEU:HG	2.10	0.81
4:T:259:ALA:O	4:T:260:ILE:HG13	1.80	0.81
3:S:117:ARG:NE	3:S:191:ARG:NH1	2.29	0.81
4:T:46:ILE:CG2	4:T:260:ILE:HD11	2.10	0.81
3:S:9:LEU:CD2	3:S:9:LEU:C	2.48	0.80
3:S:160:LEU:O	3:S:164:LEU:HD22	1.81	0.80
3:S:160:LEU:CB	3:S:163:VAL:CG1	2.55	0.80
3:S:160:LEU:CB	3:S:163:VAL:HG11	2.09	0.80
4:T:62:LEU:O	4:T:66:LEU:CG	2.30	0.80
4:T:82:LEU:HG	4:T:86:PHE:HE2	1.43	0.80
3:S:3:ARG:O	3:S:6:THR:CG2	2.30	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:196:ARG:O	3:S:199:PRO:CG	2.30	0.80
3:S:198:LYS:NZ	3:S:198:LYS:CB	2.45	0.80
3:S:173:ILE:HD12	3:S:174:PRO:HD3	1.64	0.80
1:A:16:THR:HG23	1:A:17:PRO:HD2	1.64	0.79
4:T:62:LEU:O	4:T:66:LEU:CD2	2.30	0.79
3:S:117:ARG:NE	3:S:191:ARG:HH12	1.80	0.79
3:S:185:LEU:CD1	3:S:185:LEU:O	2.30	0.79
4:T:233:HIS:N	4:T:233:HIS:ND1	2.30	0.79
4:T:92:HIS:ND1	4:T:94:TYR:CE2	2.51	0.79
4:T:85:LEU:CD2	4:T:85:LEU:O	2.30	0.79
3:S:197:LEU:C	3:S:199:PRO:HD2	2.03	0.79
4:T:84:LEU:O	4:T:88:PRO:CD	2.30	0.79
3:S:198:LYS:HZ2	3:S:198:LYS:HB2	1.46	0.79
4:T:60:ILE:O	4:T:60:ILE:CD1	2.30	0.79
3:S:160:LEU:HD22	3:S:163:VAL:HG11	1.65	0.78
4:T:62:LEU:O	4:T:66:LEU:CD1	2.30	0.78
3:S:198:LYS:N	3:S:199:PRO:CD	2.46	0.78
3:S:134:LEU:CD2	3:S:134:LEU:N	2.45	0.78
1:A:173:THR:HB	1:A:181:ARG:HG3	1.65	0.78
3:S:196:ARG:O	3:S:199:PRO:CD	2.31	0.78
4:T:83:GLN:HE22	4:T:113:ILE:HG13	1.49	0.78
3:S:185:LEU:HD22	4:T:202:ILE:HG12	1.66	0.77
4:T:40:ASN:HB2	4:T:43:SER:HB3	1.66	0.77
3:S:160:LEU:O	3:S:163:VAL:CG1	2.32	0.77
3:S:185:LEU:O	3:S:189:PRO:HD2	1.84	0.77
4:T:92:HIS:CB	4:T:94:TYR:CD2	2.65	0.75
3:S:160:LEU:O	3:S:164:LEU:HD23	1.84	0.75
3:S:185:LEU:C	3:S:185:LEU:HD13	2.07	0.75
3:S:195:GLU:O	3:S:199:PRO:CD	2.30	0.75
2:B:117:MET:HE1	4:T:235:SER:HB2	1.69	0.75
3:S:143:TYR:CD1	3:S:161:GLY:HA2	2.22	0.74
3:S:95:ARG:N	3:S:95:ARG:HD2	2.02	0.74
2:B:97:ALA:O	2:B:138:PRO:CD	2.36	0.74
2:B:46:LYS:HD2	2:B:197:ILE:HG22	1.69	0.74
4:T:85:LEU:C	4:T:85:LEU:HD23	2.07	0.74
4:T:77:VAL:O	4:T:81:VAL:CG1	2.36	0.74
2:B:201:ILE:HD13	2:B:239:LEU:CD2	2.18	0.74
4:T:46:ILE:HD13	4:T:260:ILE:HG12	1.69	0.74
3:S:160:LEU:CD2	3:S:163:VAL:HG11	2.17	0.73
4:T:87:SER:HB2	4:T:88:PRO:HD3	1.69	0.73
3:S:47:VAL:CG2	3:S:190:LEU:HD22	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:TYR:HA	1:A:141:ARG:HD3	1.70	0.73
1:A:178:PRO:HD3	2:B:199:HIS:CB	2.19	0.73
3:S:33:PHE:H	3:S:34:GLY:HA2	1.54	0.73
3:S:166:ILE:CG1	4:T:86:PHE:CZ	2.71	0.72
2:B:111:GLN:HE22	4:T:233:HIS:CD2	2.03	0.72
1:A:18:MET:O	1:A:19:ALA:O	2.07	0.72
3:S:198:LYS:N	3:S:199:PRO:HD2	2.05	0.72
2:B:190:GLN:O	2:B:190:GLN:HG3	1.89	0.72
3:S:143:TYR:HA	3:S:164:LEU:HG	0.78	0.72
3:S:139:VAL:CG1	3:S:168:LEU:CD2	2.67	0.72
4:T:51:THR:CG2	4:T:52:VAL:N	2.53	0.72
2:B:36:LEU:HD21	2:B:211:LEU:HD22	1.70	0.71
2:B:21:PRO:HB2	2:B:45:GLY:HA2	1.71	0.71
4:T:83:GLN:NE2	4:T:113:ILE:HG13	2.05	0.70
3:S:196:ARG:C	3:S:199:PRO:HD2	2.11	0.70
4:T:256:GLY:O	4:T:260:ILE:CG1	2.39	0.70
3:S:185:LEU:CD2	4:T:202:ILE:CG1	2.66	0.70
1:A:171:GLU:HA	1:A:171:GLU:OE2	1.90	0.70
1:A:100:GLU:HB2	1:A:106:ASP:HB2	1.73	0.69
3:S:182:VAL:O	3:S:186:ILE:HG22	1.93	0.69
4:T:228:TYR:O	4:T:228:TYR:HD1	1.76	0.69
1:A:256:ARG:NE	2:B:269:GLU:OE1	2.25	0.69
3:S:47:VAL:CA	3:S:190:LEU:HD22	2.18	0.68
3:S:197:LEU:HD22	3:S:197:LEU:O	1.93	0.68
2:B:114:ARG:HG2	4:T:236:GLN:HA	1.74	0.68
4:T:51:THR:O	4:T:55:ILE:HG12	1.94	0.68
4:T:46:ILE:HG21	4:T:260:ILE:HD11	1.68	0.68
4:T:23:ALA:HB1	4:T:250:TRP:HE1	1.58	0.68
3:S:198:LYS:HZ2	3:S:198:LYS:CB	2.06	0.68
1:A:153:ARG:NH2	1:A:183:GLU:OE1	2.27	0.67
3:S:47:VAL:CA	3:S:190:LEU:HD21	2.17	0.67
3:S:139:VAL:HG12	3:S:168:LEU:CD2	2.22	0.67
4:T:43:SER:OG	4:T:44:TYR:N	2.28	0.67
3:S:196:ARG:C	3:S:199:PRO:CD	2.63	0.67
4:T:228:TYR:C	4:T:228:TYR:CD1	2.67	0.66
4:T:228:TYR:O	4:T:228:TYR:CD1	2.47	0.66
4:T:106:GLY:O	4:T:110:ALA:HB3	1.96	0.66
4:T:60:ILE:CG1	4:T:60:ILE:O	2.43	0.66
4:T:65:PHE:CZ	4:T:122:MET:HA	2.30	0.66
4:T:63:GLY:O	4:T:66:LEU:CD1	2.30	0.66
4:T:128:THR:HG23	4:T:129:LEU:HD13	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:166:ILE:HG12	4:T:86:PHE:CZ	2.31	0.65
4:T:64:PHE:O	4:T:68:GLY:N	2.28	0.65
2:B:172:ASP:OD1	2:B:172:ASP:N	2.30	0.65
3:S:47:VAL:CB	3:S:190:LEU:HD22	2.27	0.65
1:A:173:THR:CB	1:A:181:ARG:HG3	2.26	0.65
1:A:178:PRO:HD3	2:B:199:HIS:HB2	1.78	0.65
3:S:163:VAL:HG13	3:S:164:LEU:HD22	1.79	0.65
2:B:98:THR:HG22	2:B:101:ASP:OD2	1.97	0.65
4:T:257:PHE:HA	4:T:260:ILE:HD12	1.79	0.64
4:T:62:LEU:C	4:T:66:LEU:HD11	2.17	0.64
4:T:65:PHE:CD2	4:T:69:ILE:CD1	2.81	0.64
3:S:200:GLN:N	3:S:200:GLN:OE1	2.30	0.64
1:A:282:TRP:O	1:A:285:ARG:HB2	1.96	0.64
3:S:196:ARG:O	3:S:200:GLN:NE2	2.30	0.64
2:B:10:HIS:HB2	2:B:61:GLN:HB2	1.78	0.64
4:T:87:SER:HB2	4:T:88:PRO:CD	2.27	0.64
1:A:284:GLN:O	1:A:285:ARG:C	2.34	0.64
3:S:143:TYR:HB2	3:S:164:LEU:CG	2.28	0.64
3:S:166:ILE:HG13	4:T:86:PHE:CZ	2.32	0.64
1:A:54:ASN:HB3	1:A:56:LEU:HD22	1.80	0.64
3:S:85:PHE:CD1	3:S:91:SER:HB2	2.32	0.63
4:T:38:ALA:HB1	4:T:44:TYR:HB3	1.81	0.63
3:S:160:LEU:HB3	3:S:163:VAL:HG13	1.75	0.63
1:A:169:LEU:HB2	1:A:172:PRO:HG3	1.81	0.63
4:T:77:VAL:O	4:T:81:VAL:HG13	1.96	0.63
3:S:198:LYS:HB3	3:S:198:LYS:NZ	2.14	0.63
3:S:95:ARG:HD2	3:S:95:ARG:H	1.62	0.63
3:S:46:ILE:CB	3:S:186:ILE:HD12	2.22	0.63
4:T:92:HIS:HB3	4:T:94:TYR:CE2	2.32	0.63
4:T:51:THR:HG23	4:T:52:VAL:N	2.14	0.62
3:S:198:LYS:HB3	3:S:198:LYS:HZ3	1.63	0.62
2:B:65:THR:HA	2:B:70:THR:HA	1.81	0.62
2:B:201:ILE:CD1	2:B:239:LEU:CD2	2.78	0.62
4:T:182:GLN:O	4:T:187:VAL:HG23	1.99	0.62
4:T:106:GLY:O	4:T:110:ALA:CB	2.48	0.62
2:B:117:MET:HG2	4:T:236:GLN:HG3	1.81	0.62
2:B:11:LEU:HD12	2:B:59:PHE:HB3	1.81	0.62
1:A:92:GLN:HG3	1:A:171:GLU:HB2	1.82	0.62
4:T:46:ILE:HG22	4:T:260:ILE:CD1	2.29	0.61
2:B:201:ILE:HD13	2:B:239:LEU:HG	1.83	0.61
3:S:33:PHE:N	3:S:34:GLY:HA2	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:65:PHE:HZ	4:T:122:MET:HA	1.64	0.61
1:A:190:ARG:O	1:A:194:GLU:CG	2.47	0.61
2:B:3:GLU:N	2:B:160:LYS:HZ1	1.98	0.61
4:T:46:ILE:CG2	4:T:260:ILE:HG12	2.15	0.61
3:S:36:PHE:HB3	3:S:174:PRO:HG3	1.81	0.61
3:S:182:VAL:O	3:S:186:ILE:HB	2.01	0.61
4:T:77:VAL:O	4:T:81:VAL:HG12	1.99	0.61
4:T:52:VAL:HA	4:T:55:ILE:HG13	1.83	0.61
2:B:231:GLY:O	2:B:234:LEU:HB2	2.01	0.61
3:S:188:MET:O	3:S:192:ARG:NE	2.34	0.60
3:S:91:SER:O	3:S:95:ARG:HD3	2.01	0.60
3:S:45:VAL:HG12	3:S:60:ILE:HG22	1.83	0.60
4:T:65:PHE:CE2	4:T:69:ILE:HD12	2.35	0.60
3:S:182:VAL:O	3:S:186:ILE:CB	2.49	0.60
4:T:146:PRO:O	4:T:148:ARG:N	2.30	0.60
4:T:67:LYS:N	4:T:67:LYS:HD2	2.17	0.60
1:A:171:GLU:OE2	1:A:171:GLU:CA	2.49	0.60
2:B:35:TRP:HB2	2:B:187:ARG:NH1	2.17	0.60
4:T:48:ILE:O	4:T:51:THR:HG22	2.02	0.60
3:S:200:GLN:CA	3:S:200:GLN:OE1	2.50	0.59
4:T:19:LEU:N	4:T:19:LEU:HD12	2.16	0.59
1:A:135:ASP:OD2	1:A:136:GLN:N	2.35	0.59
2:B:189:GLN:O	2:B:190:GLN:HB3	2.03	0.59
3:S:139:VAL:CG1	3:S:168:LEU:HD21	2.27	0.59
2:B:109:ASN:O	4:T:227:GLY:O	2.20	0.59
3:S:43:LEU:HB2	3:S:182:VAL:HG12	1.84	0.59
3:S:185:LEU:HD22	4:T:202:ILE:CG1	2.33	0.59
4:T:259:ALA:C	4:T:260:ILE:HG13	2.23	0.58
4:T:164:ALA:O	4:T:168:VAL:HG23	2.03	0.58
3:S:182:VAL:O	3:S:186:ILE:CG2	2.50	0.58
1:A:262:LEU:HD23	1:A:281:GLN:OE1	2.02	0.58
1:A:190:ARG:O	1:A:194:GLU:HG3	2.03	0.58
1:A:172:PRO:HB3	1:A:188:PHE:HE2	1.68	0.58
4:T:92:HIS:CB	4:T:94:TYR:CE2	2.86	0.58
1:A:9:THR:HB	1:A:24:THR:HA	1.84	0.58
4:T:21:PRO:HB2	4:T:22:ARG:HG3	1.86	0.58
4:T:88:PRO:CG	4:T:103:THR:CB	2.56	0.58
4:T:93:THR:O	4:T:93:THR:HG22	2.04	0.58
2:B:230:ARG:HB2	2:B:234:LEU:HD13	1.84	0.58
4:T:132:GLN:H	4:T:132:GLN:CD	2.06	0.58
2:B:92:ASN:O	4:T:219:LEU:HA	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:VAL:HG22	1:A:166:VAL:HB	1.86	0.58
3:S:51:ALA:CB	3:S:190:LEU:CD1	2.60	0.57
3:S:9:LEU:CD2	3:S:9:LEU:O	2.30	0.57
4:T:250:TRP:O	4:T:254:LEU:HB2	2.04	0.57
2:B:105:PHE:HD2	4:T:226:ARG:HD2	1.69	0.57
2:B:26:VAL:HG13	2:B:218:LEU:HB2	1.87	0.57
2:B:240:ASP:N	2:B:240:ASP:OD2	2.24	0.57
4:T:105:ASP:N	4:T:105:ASP:OD1	2.30	0.57
1:A:18:MET:O	1:A:19:ALA:C	2.42	0.57
2:B:201:ILE:HD13	2:B:239:LEU:HD21	1.86	0.57
2:B:65:THR:HB	2:B:70:THR:HG22	1.87	0.57
2:B:81:GLU:HG3	2:B:82:GLN:HG2	1.86	0.57
2:B:46:LYS:HD2	2:B:197:ILE:CG2	2.35	0.56
4:T:65:PHE:HZ	4:T:122:MET:CA	2.18	0.56
4:T:65:PHE:CD2	4:T:69:ILE:HD12	2.41	0.56
2:B:111:GLN:CD	4:T:233:HIS:HD2	2.05	0.56
3:S:7:PHE:CD1	3:S:7:PHE:C	2.78	0.56
3:S:124:ALA:CB	3:S:186:ILE:HG21	2.36	0.56
1:A:169:LEU:HD22	1:A:188:PHE:CZ	2.39	0.56
4:T:46:ILE:HG22	4:T:260:ILE:HD13	1.88	0.56
4:T:51:THR:HG22	4:T:52:VAL:H	1.71	0.56
3:S:182:VAL:C	3:S:186:ILE:HB	2.26	0.56
1:A:13:GLN:O	1:A:16:THR:HB	2.05	0.56
3:S:128:ALA:O	3:S:131:ASN:HB3	2.05	0.56
3:S:129:LEU:C	3:S:131:ASN:H	2.10	0.56
1:A:16:THR:HG23	1:A:17:PRO:CD	2.36	0.55
3:S:88:PRO:O	3:S:92:ILE:HG13	2.06	0.55
3:S:51:ALA:CA	3:S:190:LEU:HD11	2.37	0.55
4:T:88:PRO:C	4:T:95:PHE:O	2.45	0.55
2:B:201:ILE:HD13	2:B:239:LEU:CG	2.37	0.55
4:T:85:LEU:C	4:T:85:LEU:HD22	2.25	0.55
1:A:63:THR:OG1	1:A:72:THR:HG22	2.07	0.55
1:A:194:GLU:O	1:A:195:GLN:CB	2.49	0.55
3:S:143:TYR:HB2	3:S:164:LEU:HG	1.85	0.55
1:A:170:ASP:O	1:A:202:VAL:O	2.25	0.55
3:S:185:LEU:C	3:S:185:LEU:HD12	2.25	0.55
4:T:23:ALA:O	4:T:27:LEU:N	2.29	0.55
4:T:40:ASN:OD1	4:T:40:ASN:N	2.41	0.54
2:B:114:ARG:HA	2:B:117:MET:HB3	1.89	0.54
3:S:190:LEU:O	3:S:190:LEU:HD12	2.08	0.54
1:A:142:SER:HB3	1:A:145:GLU:HG3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:S:143:TYR:CE1	3:S:161:GLY:HA2	2.43	0.54
2:B:98:THR:CG2	2:B:101:ASP:OD2	2.55	0.54
2:B:36:LEU:HD11	2:B:211:LEU:HD13	1.89	0.54
4:T:46:ILE:CG2	4:T:260:ILE:HD13	2.35	0.54
4:T:22:ARG:HB3	4:T:143:LEU:HD21	1.89	0.54
1:A:282:TRP:CH2	2:B:275:SER:HB2	2.43	0.54
2:B:95:VAL:HG23	4:T:226:ARG:NH2	2.24	0.54
4:T:63:GLY:N	4:T:66:LEU:HD11	2.21	0.53
4:T:88:PRO:O	4:T:95:PHE:O	2.26	0.53
1:A:91:PHE:CE2	4:T:181:ALA:HA	2.43	0.53
3:S:143:TYR:CB	3:S:164:LEU:CG	2.83	0.53
3:S:143:TYR:HD1	3:S:143:TYR:C	2.11	0.53
4:T:237:TYR:O	4:T:238:ARG:C	2.46	0.53
3:S:68:THR:HG21	3:S:95:ARG:HH12	1.72	0.53
3:S:40:LEU:HD23	3:S:181:LEU:HD11	1.90	0.53
3:S:170:THR:HA	3:S:173:ILE:HD11	1.91	0.53
3:S:166:ILE:HG21	4:T:86:PHE:CE2	2.44	0.53
3:S:93:LEU:HB2	3:S:94:PRO:HD3	1.90	0.53
3:S:185:LEU:HD22	4:T:202:ILE:CD1	2.39	0.53
3:S:164:LEU:O	3:S:168:LEU:HG	2.09	0.52
3:S:44:THR:HB	4:T:209:PHE:HZ	1.74	0.52
4:T:18:ARG:HB3	4:T:130:SER:CB	2.13	0.52
4:T:60:ILE:HG13	4:T:60:ILE:O	2.08	0.52
1:A:33:ARG:O	1:A:192:HIS:HE1	1.91	0.52
3:S:61:GLY:O	3:S:95:ARG:HB3	2.10	0.52
3:S:34:GLY:H	3:S:35:PRO:HD2	1.73	0.52
4:T:80:VAL:HG21	4:T:114:PHE:HB2	1.92	0.52
3:S:85:PHE:HD1	3:S:91:SER:CB	2.22	0.52
3:S:3:ARG:HA	3:S:6:THR:CG2	2.39	0.52
4:T:211:SER:O	4:T:215:ARG:HB2	2.10	0.52
3:S:143:TYR:HD1	3:S:143:TYR:O	1.92	0.52
3:S:88:PRO:O	3:S:92:ILE:HB	2.10	0.52
1:A:187:LEU:O	1:A:191:LEU:HG	2.10	0.52
3:S:85:PHE:HD1	3:S:91:SER:HB2	1.73	0.52
4:T:51:THR:CG2	4:T:52:VAL:H	2.22	0.52
4:T:259:ALA:O	4:T:260:ILE:CG1	2.56	0.52
4:T:48:ILE:HA	4:T:51:THR:HG22	1.92	0.52
1:A:190:ARG:O	1:A:194:GLU:HG2	2.10	0.51
3:S:47:VAL:HG22	3:S:190:LEU:CD2	2.31	0.51
3:S:198:LYS:HB2	3:S:198:LYS:NZ	2.15	0.51
4:T:65:PHE:O	4:T:65:PHE:HD2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:105:PHE:CD2	4:T:226:ARG:HD2	2.45	0.51
3:S:132:THR:OG1	3:S:175:GLU:HB3	2.10	0.51
3:S:90:ILE:HG12	3:S:138:LEU:HB3	1.92	0.51
3:S:166:ILE:CG1	4:T:86:PHE:CE1	2.87	0.51
4:T:88:PRO:HD3	4:T:103:THR:HA	1.93	0.50
3:S:185:LEU:O	3:S:189:PRO:CD	2.58	0.50
4:T:257:PHE:O	4:T:260:ILE:HB	2.12	0.50
3:S:124:ALA:HB1	3:S:186:ILE:HG21	1.94	0.50
3:S:143:TYR:CE1	3:S:161:GLY:CA	2.94	0.50
4:T:63:GLY:C	4:T:66:LEU:CD1	2.55	0.50
4:T:18:ARG:O	4:T:20:ASP:N	2.39	0.50
1:A:256:ARG:NH2	2:B:269:GLU:OE1	2.42	0.50
4:T:27:LEU:O	4:T:31:TYR:N	2.44	0.50
3:S:64:TRP:CE3	3:S:95:ARG:HG3	2.46	0.50
1:A:92:GLN:HG3	1:A:171:GLU:CB	2.41	0.50
4:T:237:TYR:O	4:T:238:ARG:O	2.30	0.50
4:T:65:PHE:O	4:T:69:ILE:HG13	2.11	0.50
2:B:107:LEU:O	2:B:110:ARG:N	2.44	0.50
2:B:137:GLU:HG3	2:B:139:SER:H	1.77	0.50
1:A:206:MET:HE1	1:A:245:LEU:HD22	1.92	0.50
1:A:191:LEU:O	1:A:194:GLU:O	2.30	0.49
1:A:132:VAL:HG13	1:A:153:ARG:HB3	1.94	0.49
3:S:143:TYR:CD1	3:S:143:TYR:O	2.64	0.49
2:B:198:THR:HG1	2:B:200:ASP:H	1.55	0.49
4:T:142:SER:O	4:T:145:LYS:HB3	2.12	0.49
3:S:182:VAL:HG12	3:S:186:ILE:HG13	1.94	0.49
2:B:8:VAL:H	2:B:27:SER:HG	1.57	0.49
2:B:249:ALA:O	2:B:253:GLN:HG3	2.12	0.49
2:B:189:GLN:O	2:B:190:GLN:CB	2.60	0.49
1:A:134:LEU:HD21	1:A:153:ARG:HD2	1.94	0.49
4:T:48:ILE:HA	4:T:51:THR:CG2	2.42	0.49
3:S:135:VAL:O	3:S:139:VAL:HG23	2.13	0.49
4:T:125:THR:HA	4:T:128:THR:HG22	1.94	0.49
4:T:144:MET:HE3	4:T:154:VAL:HG23	1.95	0.49
1:A:238:GLU:H	1:A:238:GLU:CD	2.16	0.49
2:B:87:PHE:HB2	2:B:93:GLN:OE1	2.12	0.49
3:S:38:MET:SD	3:S:177:ILE:HG22	2.52	0.48
1:A:81:LYS:HZ2	1:A:84:ARG:HG2	1.78	0.48
3:S:192:ARG:HD3	3:S:192:ARG:H	1.78	0.48
1:A:173:THR:CG2	1:A:181:ARG:HG3	2.43	0.48
4:T:235:SER:O	4:T:236:GLN:CB	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:80:VAL:HG13	4:T:110:ALA:HB1	1.96	0.48
4:T:84:LEU:O	4:T:88:PRO:HG2	2.14	0.48
3:S:199:PRO:HG2	3:S:200:GLN:NE2	2.28	0.48
4:T:193:GLY:HA2	4:T:197:GLN:HB2	1.94	0.48
2:B:97:ALA:O	2:B:138:PRO:HD2	2.13	0.48
4:T:20:ASP:OD2	4:T:20:ASP:N	2.47	0.48
3:S:90:ILE:HA	3:S:138:LEU:HD12	1.95	0.48
4:T:46:ILE:CD1	4:T:260:ILE:HG12	2.41	0.48
3:S:131:ASN:O	3:S:135:VAL:HG22	2.13	0.48
3:S:24:LEU:HD23	3:S:24:LEU:HA	1.63	0.48
1:A:41:HIS:CE1	1:A:44:SER:HB3	2.49	0.48
2:B:201:ILE:CD1	2:B:239:LEU:HD23	2.45	0.47
2:B:55:GLY:H	2:B:80:ARG:HE	1.61	0.47
2:B:201:ILE:HD11	2:B:239:LEU:HD23	1.97	0.47
3:S:166:ILE:HA	3:S:166:ILE:HD12	1.66	0.47
1:A:91:PHE:CD2	4:T:181:ALA:HA	2.50	0.47
3:S:73:PHE:CZ	3:S:86:THR:HG22	2.49	0.47
2:B:268:MET:HB2	2:B:268:MET:HE2	1.54	0.47
3:S:117:ARG:HG3	3:S:191:ARG:HH11	1.79	0.47
3:S:200:GLN:OE1	3:S:200:GLN:HA	2.15	0.47
2:B:272:LEU:HA	2:B:275:SER:OG	2.15	0.47
3:S:22:GLN:OE1	3:S:44:THR:HG21	2.14	0.47
2:B:186:LEU:O	2:B:189:GLN:O	2.33	0.47
4:T:147:LEU:O	4:T:147:LEU:HD23	2.15	0.47
1:A:252:GLN:O	1:A:255:ARG:HB3	2.15	0.47
4:T:62:LEU:HD22	4:T:62:LEU:HA	1.71	0.46
2:B:151:LEU:O	2:B:154:ILE:HG13	2.15	0.46
1:A:178:PRO:HD3	2:B:199:HIS:HB3	1.96	0.46
3:S:98:MET:HB2	3:S:127:ALA:O	2.16	0.46
1:A:50:ILE:HG13	1:A:168:VAL:CG1	2.45	0.46
1:A:94:PRO:O	1:A:97:GLN:HG2	2.15	0.46
4:T:62:LEU:O	4:T:66:LEU:HG	2.15	0.46
4:T:84:LEU:O	4:T:88:PRO:CG	2.63	0.46
4:T:232:GLU:HB3	4:T:233:HIS:CE1	2.49	0.46
1:A:16:THR:HG22	1:A:18:MET:H	1.80	0.46
2:B:108:GLU:HB2	4:T:235:SER:O	2.15	0.46
1:A:262:LEU:CD2	1:A:281:GLN:OE1	2.63	0.46
1:A:21:THR:HG22	1:A:24:THR:HG22	1.96	0.46
2:B:95:VAL:HG23	4:T:226:ARG:HH21	1.81	0.46
1:A:52:GLN:HG3	1:A:64:ILE:HD11	1.97	0.46
1:A:239:TRP:O	1:A:243:HIS:ND1	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:174:GLU:O	4:T:178:ILE:HG13	2.16	0.46
3:S:160:LEU:O	3:S:163:VAL:HG12	2.11	0.46
1:A:170:ASP:HA	1:A:202:VAL:HG23	1.98	0.46
4:T:138:THR:HG21	4:T:237:TYR:CE1	2.51	0.46
3:S:143:TYR:CA	3:S:164:LEU:CG	2.55	0.46
3:S:34:GLY:N	3:S:35:PRO:HD2	2.30	0.46
2:B:248:LYS:O	2:B:252:ARG:HG3	2.16	0.46
4:T:86:PHE:O	4:T:90:GLY:N	2.43	0.46
4:T:20:ASP:HB2	4:T:24:LYS:HG3	1.98	0.45
4:T:81:VAL:HG22	4:T:82:LEU:N	2.31	0.45
1:A:147:SER:HB3	1:A:150:GLN:CD	2.37	0.45
2:B:12:THR:C	2:B:14:GLN:H	2.20	0.45
2:B:264:THR:H	2:B:267:GLU:HB2	1.81	0.45
1:A:207:GLU:O	1:A:211:GLN:HB3	2.17	0.45
1:A:8:VAL:HG13	1:A:59:PRO:HB3	1.98	0.45
2:B:91:ASP:OD1	2:B:91:ASP:N	2.47	0.45
3:S:14:LEU:HD11	4:T:167:PHE:CE1	2.51	0.45
3:S:39:THR:N	3:S:178:LEU:HD21	2.32	0.45
4:T:65:PHE:C	4:T:65:PHE:CD2	2.90	0.45
4:T:65:PHE:CZ	4:T:122:MET:HG2	2.52	0.45
4:T:89:ALA:O	4:T:94:TYR:HB2	2.16	0.45
4:T:18:ARG:HB2	4:T:130:SER:O	2.17	0.45
3:S:43:LEU:HD11	3:S:185:LEU:HD12	1.99	0.45
4:T:18:ARG:N	4:T:19:LEU:HD13	2.32	0.45
2:B:9:ASP:HA	2:B:62:GLY:HA3	1.97	0.45
1:A:118:ASP:N	1:A:118:ASP:OD2	2.42	0.45
3:S:47:VAL:HG13	3:S:190:LEU:HD22	1.97	0.45
3:S:117:ARG:HG3	3:S:191:ARG:NH1	2.32	0.45
3:S:116:MET:O	3:S:120:MET:HG2	2.17	0.45
4:T:84:LEU:HG	4:T:106:GLY:C	2.38	0.45
3:S:94:PRO:HG3	3:S:134:LEU:HB2	1.99	0.45
1:A:94:PRO:HB2	1:A:155:ALA:HB1	2.00	0.44
3:S:16:MET:HE3	4:T:160:MET:HB3	1.99	0.44
1:A:28:LEU:HD12	1:A:28:LEU:O	2.17	0.44
4:T:39:ASN:H	4:T:39:ASN:ND2	2.15	0.44
2:B:35:TRP:N	2:B:187:ARG:HH12	2.15	0.44
2:B:105:PHE:CZ	4:T:223:MET:HE3	2.52	0.44
3:S:76:PRO:HD2	4:T:72:LEU:HD22	1.99	0.44
1:A:278:LEU:HD23	1:A:278:LEU:HA	1.78	0.44
4:T:256:GLY:O	4:T:260:ILE:HG13	2.18	0.44
4:T:18:ARG:HD2	4:T:130:SER:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:6:ILE:HD13	2:B:28:PHE:CZ	2.53	0.44
1:A:276:ASP:OD1	2:B:254:ARG:HD3	2.17	0.44
4:T:106:GLY:O	4:T:110:ALA:N	2.45	0.44
3:S:129:LEU:O	3:S:131:ASN:N	2.49	0.44
3:S:188:MET:HB2	3:S:189:PRO:HD3	2.00	0.44
3:S:171:ASN:C	3:S:173:ILE:H	2.21	0.44
4:T:196:LYS:HB2	4:T:196:LYS:HE2	1.87	0.44
2:B:114:ARG:HG2	4:T:236:GLN:CA	2.45	0.44
4:T:132:GLN:HA	4:T:133:PRO:HD3	1.89	0.44
2:B:138:PRO:HG3	4:T:166:ARG:HE	1.83	0.43
1:A:277:TYR:O	1:A:280:GLN:HG3	2.18	0.43
3:S:139:VAL:CG1	3:S:168:LEU:HD23	2.45	0.43
4:T:65:PHE:C	4:T:65:PHE:HD2	2.20	0.43
3:S:192:ARG:H	3:S:192:ARG:CD	2.28	0.43
2:B:3:GLU:HA	2:B:30:VAL:O	2.17	0.43
3:S:188:MET:H	3:S:188:MET:HG2	1.54	0.43
1:A:18:MET:H	1:A:18:MET:HG2	1.44	0.43
4:T:235:SER:O	4:T:236:GLN:HB2	2.17	0.43
2:B:274:GLN:HG2	2:B:275:SER:N	2.34	0.43
1:A:81:LYS:HD2	1:A:81:LYS:HA	1.70	0.43
3:S:143:TYR:CB	3:S:164:LEU:CB	2.81	0.43
4:T:182:GLN:HG3	4:T:187:VAL:HG21	2.00	0.43
3:S:136:LEU:HD23	3:S:136:LEU:N	2.33	0.43
3:S:182:VAL:CB	3:S:186:ILE:HB	2.46	0.43
2:B:98:THR:CG2	2:B:101:ASP:CG	2.88	0.43
2:B:243:PHE:HA	2:B:246:LYS:HG3	2.00	0.43
4:T:132:GLN:HG2	4:T:135:ASP:HB2	2.00	0.42
2:B:177:ILE:HD12	2:B:177:ILE:HA	1.81	0.42
3:S:19:VAL:O	3:S:23:ASN:ND2	2.48	0.42
2:B:201:ILE:HG22	2:B:202:ASP:N	2.34	0.42
1:A:75:THR:HG21	1:A:80:LEU:HD21	2.01	0.42
1:A:237:ARG:HG3	1:A:247:VAL:HG13	2.02	0.42
3:S:166:ILE:HG22	3:S:167:SER:N	2.34	0.42
1:A:162:MET:HE2	1:A:162:MET:HB2	1.76	0.42
3:S:200:GLN:N	3:S:200:GLN:CD	2.72	0.42
4:T:138:THR:HG21	4:T:237:TYR:HE1	1.85	0.42
2:B:154:ILE:O	2:B:159:PRO:HD3	2.19	0.42
2:B:69:ILE:HD13	2:B:78:GLN:HB3	2.00	0.42
4:T:87:SER:CB	4:T:88:PRO:HD3	2.43	0.42
4:T:28:SER:HA	4:T:31:TYR:HB3	2.02	0.42
2:B:57:LEU:HA	2:B:58:PRO:HD2	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:THR:HG22	1:A:18:MET:N	2.35	0.42
2:B:111:GLN:CD	4:T:233:HIS:CD2	2.88	0.42
2:B:198:THR:OG1	2:B:200:ASP:N	2.30	0.42
3:S:116:MET:CE	3:S:194:TRP:HB3	2.50	0.42
3:S:192:ARG:O	3:S:196:ARG:HG2	2.20	0.42
1:A:38:ILE:HB	1:A:202:VAL:HG13	2.01	0.42
3:S:15:LEU:HD13	3:S:48:ALA:HB2	2.01	0.42
4:T:87:SER:CB	4:T:88:PRO:CD	2.92	0.41
1:A:178:PRO:CD	2:B:199:HIS:HB2	2.47	0.41
2:B:245:GLU:OE2	2:B:248:LYS:NZ	2.42	0.41
2:B:48:THR:HB	2:B:57:LEU:HD21	2.02	0.41
3:S:166:ILE:HG21	4:T:86:PHE:CZ	2.55	0.41
3:S:31:ILE:HG23	3:S:38:MET:HG3	2.01	0.41
4:T:51:THR:O	4:T:55:ILE:CG1	2.64	0.41
4:T:146:PRO:HB2	4:T:147:LEU:H	1.59	0.41
3:S:63:PHE:O	3:S:67:ILE:HG13	2.20	0.41
1:A:127:GLU:O	1:A:127:GLU:HG2	2.20	0.41
2:B:256:ILE:O	2:B:258:PRO:HD3	2.20	0.41
2:B:128:VAL:HG21	2:B:152:ALA:HB2	2.02	0.41
3:S:57:GLY:O	3:S:99:GLY:HA3	2.21	0.41
1:A:81:LYS:NZ	1:A:84:ARG:HG2	2.35	0.41
2:B:16:ASP:HB3	2:B:17:GLU:H	1.66	0.41
2:B:23:LEU:HD22	2:B:26:VAL:HG21	2.03	0.41
1:A:229:THR:HG23	1:A:229:THR:H	1.63	0.41
3:S:160:LEU:C	3:S:163:VAL:HG12	2.41	0.41
1:A:16:THR:CG2	1:A:17:PRO:N	2.84	0.41
2:B:89:ASN:HA	2:B:90:PRO:HD3	1.92	0.41
3:S:160:LEU:C	3:S:164:LEU:HD23	2.41	0.41
2:B:133:PHE:O	2:B:136:ARG:HB2	2.22	0.40
4:T:63:GLY:O	4:T:67:LYS:HD3	2.21	0.40
3:S:117:ARG:O	3:S:121:GLN:HG3	2.21	0.40
3:S:143:TYR:CE1	3:S:161:GLY:HA3	2.56	0.40
2:B:107:LEU:HB2	2:B:117:MET:SD	2.61	0.40
1:A:50:ILE:HD12	1:A:50:ILE:HA	1.60	0.40
3:S:132:THR:HA	3:S:135:VAL:CG2	2.52	0.40
3:S:173:ILE:O	3:S:177:ILE:HD13	2.21	0.40
2:B:201:ILE:HG12	2:B:240:ASP:O	2.22	0.40
3:S:89:LEU:HD23	3:S:89:LEU:HA	1.89	0.40
3:S:47:VAL:CG1	3:S:190:LEU:HD22	2.51	0.40
1:A:55:ALA:HB2	1:A:83:LEU:HD23	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	282/290 (97%)	259 (92%)	18 (6%)	5 (2%)	11 50
2	B	273/279 (98%)	244 (89%)	25 (9%)	4 (2%)	13 53
3	S	183/203 (90%)	155 (85%)	21 (12%)	7 (4%)	4 26
4	T	230/280 (82%)	196 (85%)	25 (11%)	9 (4%)	4 26
All	All	968/1052 (92%)	854 (88%)	89 (9%)	25 (3%)	7 38

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	ALA
3	S	146	PRO
4	T	146	PRO
4	T	147	LEU
1	A	78	ALA
1	A	196	GLY
4	T	19	LEU
4	T	21	PRO
4	T	150	VAL
4	T	260	ILE
2	B	16	ASP
3	S	184	PRO
3	S	190	LEU
4	T	38	ALA
2	B	41	HIS
3	S	37	SER
3	S	130	THR
3	S	173	ILE
4	T	236	GLN
1	A	55	ALA
1	A	99	PHE
2	B	159	PRO

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Mol	Chain	Res	Type
4	T	43	SER
2	B	158	ALA
3	S	172	GLY

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	233/237 (98%)	199 (85%)	34 (15%)	4 18
2	B	230/234 (98%)	180 (78%)	50 (22%)	1 6
3	S	155/167 (93%)	111 (72%)	44 (28%)	0 1
4	T	191/243 (79%)	132 (69%)	59 (31%)	0 1
All	All	809/881 (92%)	622 (77%)	187 (23%)	1 4

All (187) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	11	THR
1	A	18	MET
1	A	24	THR
1	A	35	TYR
1	A	42	THR
1	A	47	SER
1	A	49	LEU
1	A	50	ILE
1	A	52	GLN
1	A	56	LEU
1	A	63	THR
1	A	65	LYS
1	A	84	ARG
1	A	89	MET
1	A	96	ASN
1	A	102	THR
1	A	104	ARG

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Mol	Chain	Res	Type
1	A	113	ASN
1	A	116	MET
1	A	118	ASP
1	A	120	ASP
1	A	127	GLU
1	A	132	VAL
1	A	147	SER
1	A	150	GLN
1	A	169	LEU
1	A	171	GLU
1	A	195	GLN
1	A	202	VAL
1	A	203	THR
1	A	214	GLU
1	A	223	ARG
1	A	262	LEU
2	B	13	TYR
2	B	19	GLN
2	B	31	HIS
2	B	47	SER
2	B	54	ASP
2	B	56	LEU
2	B	60	THR
2	B	65	THR
2	B	70	THR
2	B	71	LEU
2	B	72	THR
2	B	74	GLU
2	B	82	GLN
2	B	91	ASP
2	B	92	ASN
2	B	93	GLN
2	B	95	VAL
2	B	110	ARG
2	B	114	ARG
2	B	131	THR
2	B	136	ARG
2	B	149	VAL
2	B	169	SER
2	B	172	ASP
2	B	177	ILE
2	B	186	LEU

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Mol	Chain	Res	Type
2	B	187	ARG
2	B	189	GLN
2	B	190	GLN
2	B	197	ILE
2	B	198	THR
2	B	199	HIS
2	B	201	ILE
2	B	202	ASP
2	B	217	ARG
2	B	218	LEU
2	B	219	VAL
2	B	223	VAL
2	B	234	LEU
2	B	236	GLU
2	B	240	ASP
2	B	244	THR
2	B	245	GLU
2	B	246	LYS
2	B	247	LEU
2	B	254	ARG
2	B	260	THR
2	B	269	GLU
2	B	274	GLN
2	B	275	SER
3	S	4	HIS
3	S	5	LYS
3	S	6	THR
3	S	7	PHE
3	S	8	ARG
3	S	9	LEU
3	S	14	LEU
3	S	21	LEU
3	S	31	ILE
3	S	33	PHE
3	S	43	LEU
3	S	44	THR
3	S	45	VAL
3	S	55	ARG
3	S	59	LEU
3	S	77	SER
3	S	86	THR
3	S	91	SER

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Mol	Chain	Res	Type
3	S	92	ILE
3	S	93	LEU
3	S	95	ARG
3	S	98	MET
3	S	104	SER
3	S	107	LEU
3	S	112	ARG
3	S	131	ASN
3	S	134	LEU
3	S	135	VAL
3	S	143	TYR
3	S	144	GLN
3	S	145	THR
3	S	162	TYR
3	S	166	ILE
3	S	169	PHE
3	S	173	ILE
3	S	181	LEU
3	S	185	LEU
3	S	186	ILE
3	S	188	MET
3	S	190	LEU
3	S	192	ARG
3	S	197	LEU
3	S	198	LYS
3	S	200	GLN
4	T	18	ARG
4	T	19	LEU
4	T	20	ASP
4	T	22	ARG
4	T	26	MET
4	T	27	LEU
4	T	30	CYS
4	T	39	ASN
4	T	40	ASN
4	T	43	SER
4	T	44	TYR
4	T	50	PHE
4	T	51	THR
4	T	55	ILE
4	T	56	LEU
4	T	57	SER

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Mol	Chain	Res	Type
4	T	58	SER
4	T	60	ILE
4	T	62	LEU
4	T	65	PHE
4	T	66	LEU
4	T	70	ARG
4	T	74	TRP
4	T	78	PHE
4	T	79	THR
4	T	80	VAL
4	T	84	LEU
4	T	85	LEU
4	T	105	ASP
4	T	107	LEU
4	T	112	TYR
4	T	126	LEU
4	T	129	LEU
4	T	132	GLN
4	T	145	LYS
4	T	147	LEU
4	T	148	ARG
4	T	155	ASP
4	T	166	ARG
4	T	174	GLU
4	T	177	LYS
4	T	182	GLN
4	T	185	ARG
4	T	194	LEU
4	T	201	LEU
4	T	202	ILE
4	T	208	LEU
4	T	224	GLU
4	T	228	TYR
4	T	229	GLN
4	T	230	ASP
4	T	233	HIS
4	T	234	ARG
4	T	235	SER
4	T	236	GLN
4	T	237	TYR
4	T	247	THR
4	T	254	LEU

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Mol	Chain	Res	Type
4	T	261	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	192	HIS
2	B	109	ASN
2	B	111	GLN
2	B	274	GLN
4	T	83	GLN
4	T	132	GLN
4	T	233	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	284/290 (97%)	0.23	9 (3%) 51 39	28, 58, 99, 154	0
2	B	275/279 (98%)	0.34	14 (5%) 32 21	35, 80, 139, 274	0
3	S	187/203 (92%)	0.27	5 (2%) 58 46	44, 87, 134, 218	0
4	T	236/280 (84%)	0.52	26 (11%) 7 5	37, 103, 170, 212	0
All	All	982/1052 (93%)	0.34	54 (5%) 29 18	28, 78, 144, 274	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	18	ASN	8.7
2	B	20	ALA	8.1
4	T	234	ARG	5.1
4	T	90	GLY	5.1
2	B	14	GLN	5.0
4	T	246	ASP	4.8
2	B	19	GLN	4.7
1	A	260	ARG	4.1
1	A	21	THR	4.0
3	S	113	GLN	3.7
4	T	227	GLY	3.6
1	A	221	GLU	3.6
4	T	218	ASP	3.5
2	B	97	ALA	3.4
4	T	247	THR	3.4
3	S	115	SER	3.1
2	B	21	PRO	3.0
4	T	97	TRP	2.9
2	B	17	GLU	2.7
2	B	13	TYR	2.7
1	A	220	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	16	ASP	2.6
4	T	262	ILE	2.6
3	S	114	TRP	2.5
4	T	20	ASP	2.5
4	T	235	SER	2.5
4	T	87	SER	2.5
1	A	242	ASP	2.5
4	T	117	PHE	2.4
1	A	278	LEU	2.4
4	T	149	TRP	2.4
3	S	77	SER	2.4
2	B	89	ASN	2.4
4	T	150	VAL	2.3
3	S	118	GLN	2.3
2	B	169	SER	2.3
2	B	60	THR	2.3
4	T	256	GLY	2.2
4	T	59	LYS	2.2
4	T	65	PHE	2.2
4	T	259	ALA	2.2
4	T	86	PHE	2.2
1	A	24	THR	2.2
4	T	249	THR	2.1
4	T	60	ILE	2.1
4	T	85	LEU	2.1
1	A	22	ALA	2.1
4	T	245	ARG	2.1
2	B	96	GLY	2.1
4	T	229	GLN	2.1
4	T	88	PRO	2.1
2	B	15	TYR	2.0
1	A	60	THR	2.0
4	T	207	PRO	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 carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.