



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 01:54 am BST

PDB ID : 1SSM
Title : Serine Acetyltransferase- Apoenzyme (truncated)
Authors : Olsen, L.R.; Huang, B.; Vetting, M.W.; Roderick, S.L.
Deposited on : 2004-03-24
Resolution : 2.15 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

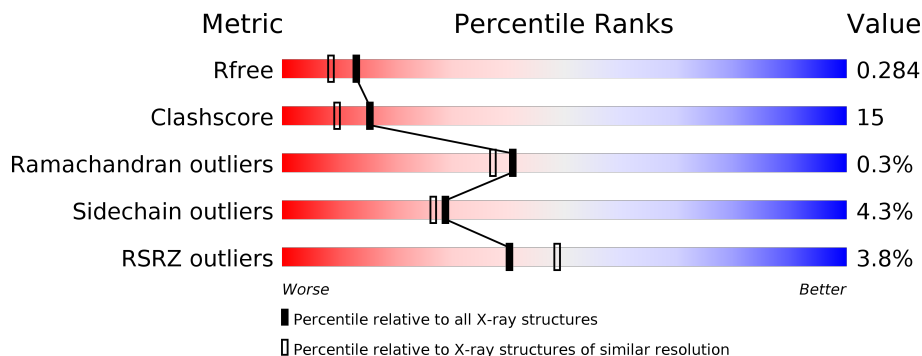
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 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	242	 5% 71% 26% ..
1	B	242	 3% 71% 26% ..
1	C	242	 3% 71% 27% ..
1	D	242	 3% 76% 22% ..
1	E	242	 5% 70% 27% ..
1	F	242	 4% 65% 31% ..

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 11514 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 Serine acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	240	1824	1162	320	336	2	4	0	0	0
1	B	240	1826	1163	319	337	2	5	0	0	0
1	C	240	1819	1159	318	336	2	4	0	0	0
1	D	240	1823	1162	319	336	2	4	0	0	0
1	E	240	1831	1166	321	337	2	5	0	0	0
1	F	240	1822	1162	319	335	2	4	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
A	22	MSE	MET	MODIFIED RESIDUE	UNP P43886
A	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
A	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
A	197	MSE	MET	MODIFIED RESIDUE	UNP P43886
B	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
B	22	MSE	MET	MODIFIED RESIDUE	UNP P43886
B	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
B	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
B	197	MSE	MET	MODIFIED RESIDUE	UNP P43886
C	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
C	22	MSE	MET	MODIFIED RESIDUE	UNP P43886
C	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
C	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
C	197	MSE	MET	MODIFIED RESIDUE	UNP P43886
D	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
D	22	MSE	MET	MODIFIED RESIDUE	UNP P43886

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Chain	Residue	Modelled	Actual	Comment	Reference
D	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
D	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
D	197	MSE	MET	MODIFIED RESIDUE	UNP P43886
E	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
E	22	MSE	MET	MODIFIED RESIDUE	UNP P43886
E	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
E	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
E	197	MSE	MET	MODIFIED RESIDUE	UNP P43886
F	1	MSE	MET	MODIFIED RESIDUE	UNP P43886
F	22	MSE	MET	MODIFIED RESIDUE	UNP P43886
F	54	MSE	MET	MODIFIED RESIDUE	UNP P43886
F	151	MSE	MET	MODIFIED RESIDUE	UNP P43886
F	197	MSE	MET	MODIFIED RESIDUE	UNP P43886

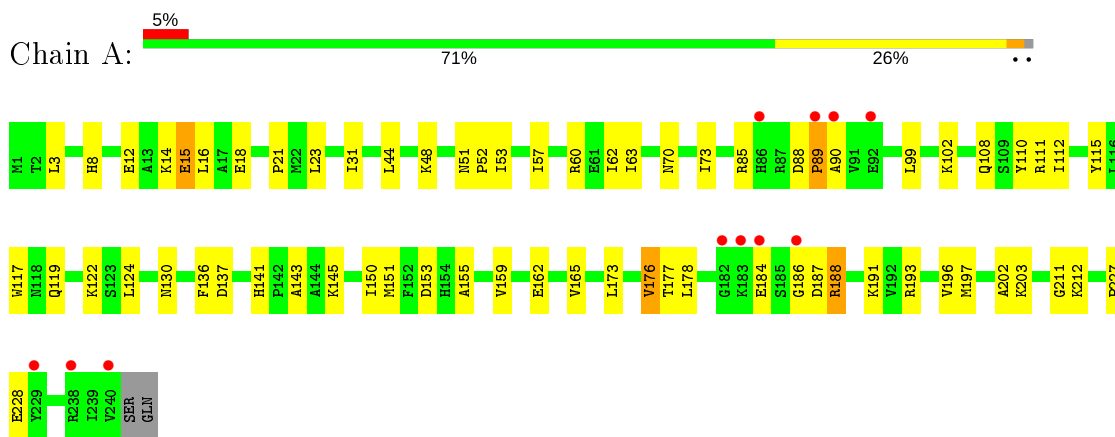
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	101	Total O 101 101	0	0
2	B	96	Total O 96 96	0	0
2	C	89	Total O 89 89	0	0
2	D	91	Total O 91 91	0	0
2	E	86	Total O 86 86	0	0
2	F	106	Total O 106 106	0	0

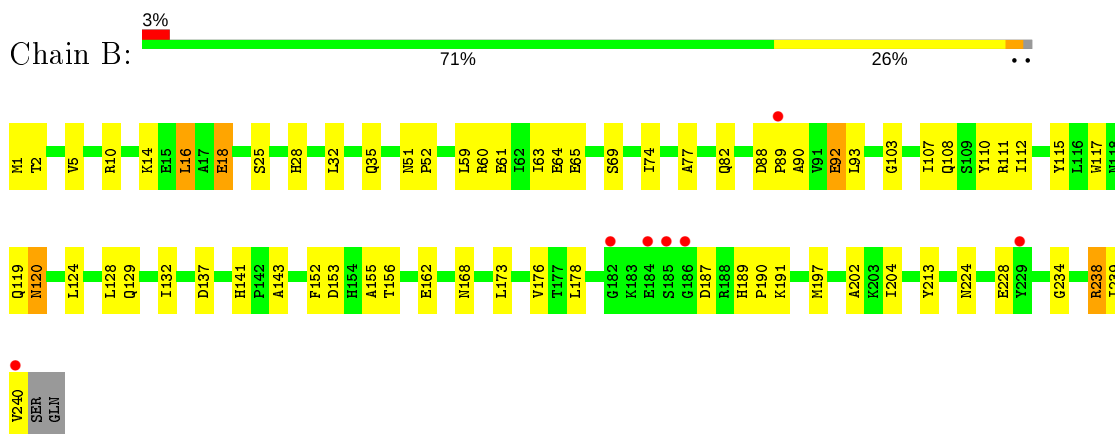
3 Residue-property plots [i](#)

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

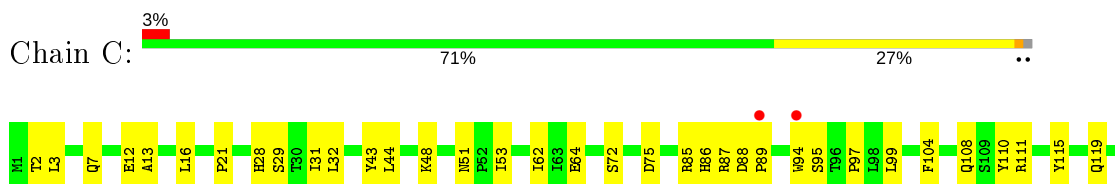
- Molecule 1: Serine acetyltransferase

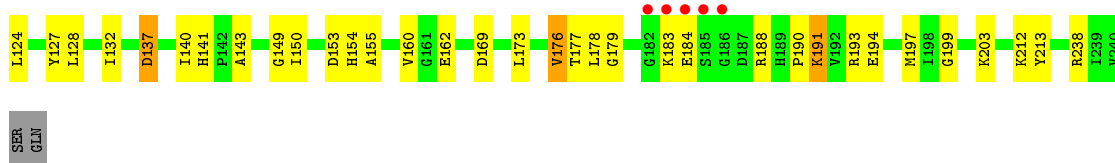


- Molecule 1: Serine acetyltransferase

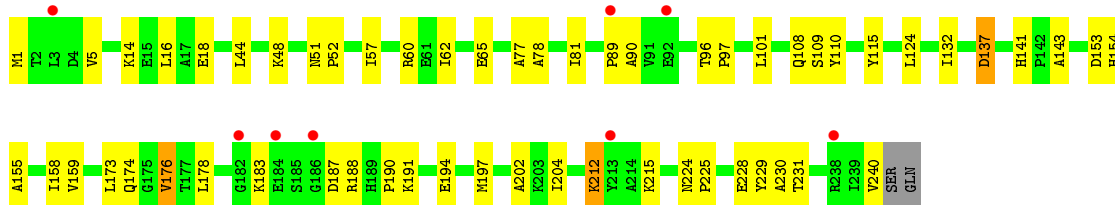
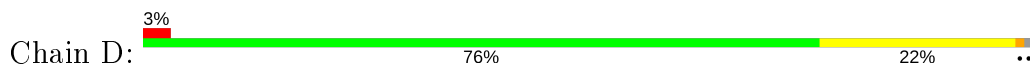


- Molecule 1: Serine acetyltransferase

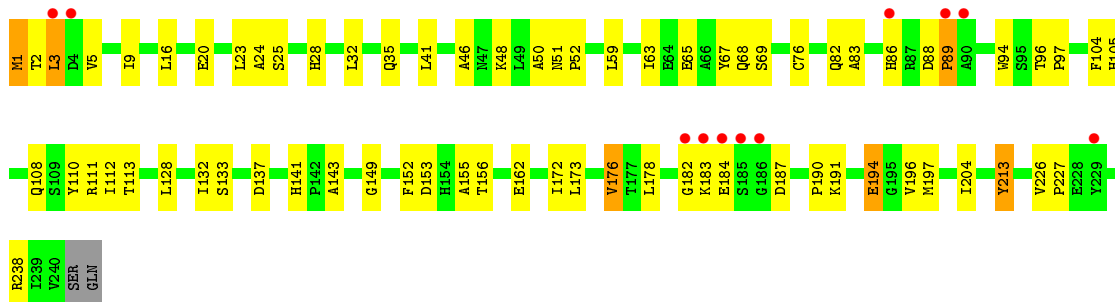




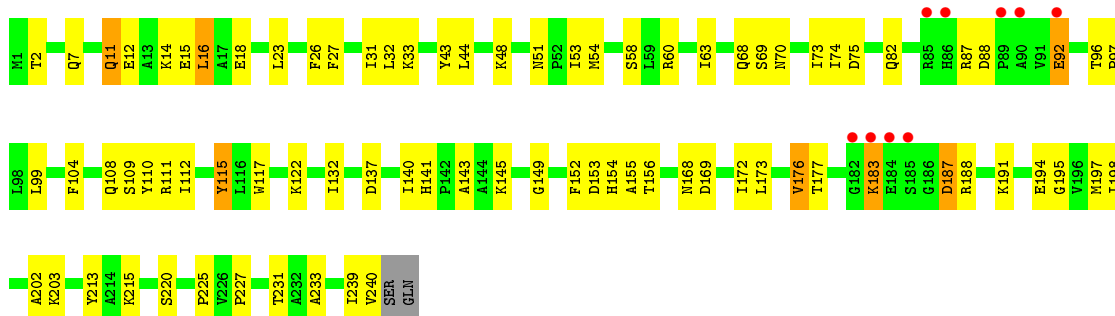
• Molecule 1: Serine acetyltransferase



• Molecule 1: Serine acetyltransferase



• Molecule 1: Serine acetyltransferase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.10Å 126.60Å 107.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 2.15 29.74 – 2.13	Depositor EDS
% Data completeness (in resolution range)	88.9 (29.74-2.15) 90.5 (29.74-2.13)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.52 (at 2.14Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.227 , 0.278 0.231 , 0.284	Depositor DCC
R_{free} test set	7181 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtrriage
Anisotropy	0.336	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.000 for l,-k,h	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11514	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.91 % 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 9.2745e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.83	2/1859 (0.1%)	0.93	3/2525 (0.1%)
1	B	0.84	1/1860 (0.1%)	0.95	2/2525 (0.1%)
1	C	0.84	0/1853	0.93	2/2517 (0.1%)
1	D	0.82	0/1857	0.93	2/2521 (0.1%)
1	E	0.85	1/1866 (0.1%)	0.93	2/2533 (0.1%)
1	F	0.81	1/1857 (0.1%)	0.92	3/2522 (0.1%)
All	All	0.83	5/11152 (0.0%)	0.93	14/15143 (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
1	B	0	1
1	C	0	1
1	E	0	1
All	All	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	76	CYS	CB-SG	7.41	1.94	1.82
1	A	15	GLU	CB-CG	5.52	1.62	1.52
1	A	15	GLU	CG-CD	5.36	1.59	1.51
1	F	43	TYR	CD2-CE2	5.22	1.47	1.39
1	B	92	GLU	CD-OE2	5.06	1.31	1.25

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	183	LYS	N-CA-C	-7.08	91.89	111.00
1	B	178	LEU	N-CA-C	-6.80	92.63	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	88	ASP	CB-CG-OD1	-6.55	112.41	118.30
1	E	3	LEU	CB-CG-CD1	-6.50	99.96	111.00
1	A	178	LEU	N-CA-C	-6.35	93.85	111.00
1	D	178	LEU	N-CA-C	-5.80	95.34	111.00
1	B	155	ALA	N-CA-C	5.49	125.82	111.00
1	A	85	ARG	NE-CZ-NH1	-5.47	117.57	120.30
1	F	60	ARG	NE-CZ-NH2	5.33	122.96	120.30
1	F	155	ALA	N-CA-C	5.33	125.38	111.00
1	F	183	LYS	N-CA-C	-5.25	96.81	111.00
1	E	178	LEU	N-CA-C	-5.24	96.86	111.00
1	A	155	ALA	N-CA-C	5.15	124.91	111.00
1	C	178	LEU	N-CA-C	-5.03	97.42	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	213	TYR	Sidechain
1	C	127	TYR	Sidechain
1	E	213	TYR	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1824	0	1821	50	0
1	B	1826	0	1829	69	0
1	C	1819	0	1816	54	0
1	D	1823	0	1827	53	0
1	E	1831	0	1834	52	0
1	F	1822	0	1823	66	0
2	A	101	0	0	14	0
2	B	96	0	0	14	0
2	C	89	0	0	11	0
2	D	91	0	0	9	0
2	E	86	0	0	4	0
2	F	106	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11514	0	10950	324	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (324) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ILE:HB	2:A:7567:HOH:O	1.47	1.12
1:E:141:HIS:HD2	1:E:143:ALA:H	1.00	0.98
1:E:3:LEU:HG	2:E:7410:HOH:O	1.65	0.95
1:E:141:HIS:CD2	1:E:143:ALA:H	1.85	0.94
1:C:141:HIS:HD2	1:C:143:ALA:H	1.04	0.94
1:B:141:HIS:HD2	1:B:143:ALA:H	1.00	0.93
1:D:141:HIS:HD2	1:D:143:ALA:H	1.11	0.91
1:B:92:GLU:HG3	1:B:93:LEU:H	1.37	0.90
1:B:141:HIS:CD2	1:B:143:ALA:H	1.90	0.89
1:F:197:MSE:HE2	1:F:215:LYS:HG2	1.54	0.89
1:D:141:HIS:CD2	1:D:143:ALA:H	1.91	0.88
1:C:141:HIS:CD2	1:C:143:ALA:H	1.92	0.87
1:C:3:LEU:O	1:C:7:GLN:HG2	1.75	0.86
1:B:238:ARG:HB3	1:B:238:ARG:NH1	1.93	0.83
1:B:92:GLU:HG3	1:B:93:LEU:HG	1.58	0.81
1:C:21:PRO:HG3	2:C:7555:HOH:O	1.80	0.81
1:C:31:ILE:HD11	1:C:99:LEU:HD21	1.61	0.80
1:B:92:GLU:HG3	1:B:93:LEU:N	1.95	0.80
1:F:7:GLN:NE2	2:F:7365:HOH:O	2.18	0.77
1:B:59:LEU:HD11	2:B:7465:HOH:O	1.85	0.76
1:E:41:LEU:HD23	1:E:67:TYR:OH	1.86	0.76
1:D:14:LYS:O	1:D:18:GLU:HG2	1.86	0.76
1:E:141:HIS:HD2	1:E:143:ALA:N	1.83	0.74
1:D:1:MSE:O	1:D:5:VAL:HG23	1.88	0.73
1:D:89:PRO:HG2	1:E:184:GLU:CB	2.19	0.73
1:F:141:HIS:HD2	1:F:143:ALA:H	1.38	0.72
1:B:187:ASP:OD1	1:B:191:LYS:NZ	2.22	0.71
1:B:88:ASP:C	1:B:88:ASP:OD1	2.28	0.71
1:A:184:GLU:CB	1:C:89:PRO:HG3	2.20	0.71
1:D:212:LYS:HG2	2:D:7323:HOH:O	1.89	0.71
1:D:89:PRO:HD2	2:D:7533:HOH:O	1.89	0.71
1:A:141:HIS:HD2	1:A:143:ALA:H	1.35	0.71
1:C:141:HIS:HD2	1:C:143:ALA:N	1.85	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:LEU:O	1:B:132:ILE:HG13	1.91	0.70
1:B:14:LYS:O	1:B:18:GLU:HG2	1.90	0.70
1:A:3:LEU:HD23	2:A:7544:HOH:O	1.92	0.70
1:F:197:MSE:HE1	1:F:215:LYS:NZ	2.06	0.70
1:F:231:THR:HG22	1:F:240:VAL:HG12	1.73	0.70
1:D:44:LEU:O	1:D:48:LYS:HG2	1.92	0.69
1:B:1:MSE:HG2	1:B:2:THR:N	2.08	0.69
1:D:197:MSE:HE1	1:D:215:LYS:HE2	1.74	0.68
1:A:187:ASP:OD1	1:A:191:LYS:NZ	2.25	0.68
1:F:153:ASP:HB3	1:F:173:LEU:HD23	1.76	0.68
1:A:141:HIS:CD2	1:A:143:ALA:H	2.10	0.68
1:E:28:HIS:NE2	2:E:7230:HOH:O	2.25	0.68
1:F:74:ILE:HD13	2:F:7261:HOH:O	1.93	0.67
1:A:14:LYS:O	1:A:18:GLU:HG2	1.94	0.67
1:D:173:LEU:O	1:D:176:VAL:HG13	1.95	0.66
1:E:1:MSE:HE3	1:E:2:THR:H	1.61	0.66
1:B:92:GLU:CG	1:B:93:LEU:H	2.07	0.66
1:A:90:ALA:N	2:A:7552:HOH:O	2.24	0.65
1:A:44:LEU:O	1:A:48:LYS:HG2	1.95	0.65
1:B:168:ASN:OD1	2:B:7374:HOH:O	2.14	0.65
1:C:85:ARG:NH1	1:C:94:TRP:CZ2	2.65	0.65
1:E:89:PRO:HA	2:E:7549:HOH:O	1.97	0.65
1:C:183:LYS:CB	2:C:7476:HOH:O	2.44	0.64
1:D:155:ALA:O	1:D:158:ILE:HG13	1.97	0.64
1:A:89:PRO:HD2	2:A:7552:HOH:O	1.97	0.64
1:C:31:ILE:CD1	1:C:99:LEU:HD21	2.27	0.64
1:F:141:HIS:CD2	1:F:143:ALA:H	2.15	0.64
1:B:137:ASP:OD1	1:B:156:THR:HG23	1.97	0.63
1:F:122:LYS:HG3	2:F:7300:HOH:O	1.97	0.63
1:D:96:THR:HB	1:D:97:PRO:HD3	1.79	0.63
1:B:1:MSE:HG2	1:B:2:THR:H	1.62	0.63
1:E:108:GLN:HE22	1:E:111:ARG:HE	1.47	0.63
1:A:62:ILE:HD13	1:A:124:LEU:HD11	1.81	0.63
1:F:197:MSE:HE1	1:F:215:LYS:HZ2	1.62	0.62
1:E:88:ASP:C	1:E:88:ASP:OD1	2.35	0.62
1:B:156:THR:HG23	1:C:137:ASP:OD2	2.00	0.62
1:C:191:LYS:HG2	1:C:193:ARG:NH2	2.14	0.62
1:A:153:ASP:HB3	1:A:173:LEU:HD23	1.81	0.62
1:B:240:VAL:O	1:B:240:VAL:HG23	2.00	0.62
1:C:173:LEU:O	1:C:176:VAL:HG13	2.01	0.61
1:B:108:GLN:HE22	1:B:111:ARG:HE	1.48	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:ARG:HB3	1:B:238:ARG:HH11	1.64	0.61
1:D:141:HIS:HD2	1:D:143:ALA:N	1.92	0.60
1:C:173:LEU:HD12	1:C:199:GLY:HA2	1.81	0.60
1:E:82:GLN:CG	1:E:94:TRP:HZ3	2.14	0.60
1:B:238:ARG:HD2	2:B:7540:HOH:O	2.02	0.60
1:F:187:ASP:OD1	1:F:191:LYS:NZ	2.29	0.60
1:B:108:GLN:HA	1:B:108:GLN:NE2	2.17	0.60
1:E:197:MSE:HG3	2:E:7493:HOH:O	2.02	0.59
1:F:197:MSE:HE2	1:F:215:LYS:CG	2.31	0.59
1:B:89:PRO:HG3	1:C:184:GLU:O	2.01	0.59
1:B:141:HIS:HD2	1:B:143:ALA:N	1.85	0.59
1:F:168:ASN:OD1	2:F:7332:HOH:O	2.16	0.59
1:C:115:TYR:O	1:C:119:GLN:HG2	2.03	0.59
1:C:128:LEU:O	1:C:132:ILE:HG13	2.01	0.59
1:A:159:VAL:HG11	1:C:154:HIS:CD2	2.38	0.59
1:F:122:LYS:HE3	2:F:7417:HOH:O	2.03	0.59
1:F:183:LYS:O	1:F:188:ARG:HD2	2.03	0.59
1:F:70:ASN:HB3	1:F:73:ILE:HD12	1.84	0.58
1:A:14:LYS:HE3	2:A:7479:HOH:O	2.04	0.58
1:F:197:MSE:CE	1:F:215:LYS:HG2	2.31	0.58
1:A:51:ASN:ND2	2:A:7165:HOH:O	2.36	0.58
1:B:153:ASP:HB3	1:B:173:LEU:HD23	1.86	0.58
1:B:238:ARG:CD	2:B:7489:HOH:O	2.52	0.57
1:B:88:ASP:OD1	1:B:90:ALA:N	2.37	0.57
1:D:65:GLU:HB3	2:D:7346:HOH:O	2.03	0.57
1:F:44:LEU:O	1:F:48:LYS:HG2	2.03	0.57
1:B:190:PRO:CG	1:B:204:ILE:HG22	2.34	0.57
1:A:70:ASN:ND2	1:A:73:ILE:HG13	2.19	0.57
1:F:109:SER:HB3	1:F:132:ILE:HD11	1.87	0.57
1:F:23:LEU:O	1:F:26:PHE:HB3	2.05	0.57
1:C:183:LYS:O	1:C:188:ARG:HD2	2.05	0.57
1:B:92:GLU:HG2	2:B:7009:HOH:O	2.04	0.56
1:E:82:GLN:HG2	1:E:94:TRP:HZ3	1.70	0.56
1:E:156:THR:CG2	1:F:137:ASP:OD2	2.53	0.56
1:B:137:ASP:OD1	1:C:137:ASP:OD2	2.23	0.56
1:D:231:THR:HG22	1:D:240:VAL:CG2	2.36	0.56
1:C:7:GLN:NE2	1:C:7:GLN:HA	2.20	0.56
1:F:183:LYS:N	2:F:7473:HOH:O	2.37	0.56
1:E:187:ASP:OD1	1:E:191:LYS:NZ	2.33	0.56
1:D:228:GLU:HG2	2:D:7338:HOH:O	2.05	0.56
1:F:88:ASP:C	1:F:88:ASP:OD1	2.43	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:ARG:HG2	2:B:7540:HOH:O	2.06	0.56
1:E:82:GLN:HG2	1:E:94:TRP:CZ3	2.40	0.56
1:F:173:LEU:O	1:F:176:VAL:HG13	2.06	0.56
1:A:23:LEU:HD22	1:B:124:LEU:HD12	1.86	0.55
1:B:65:GLU:OE1	1:B:69:SER:OG	2.21	0.55
1:B:60:ARG:O	1:B:64:GLU:HG3	2.06	0.55
1:B:28:HIS:HA	1:B:32:LEU:HB2	1.88	0.55
1:C:108:GLN:HE22	1:C:111:ARG:HE	1.55	0.55
1:B:74:ILE:O	1:B:77:ALA:HB3	2.05	0.55
1:B:238:ARG:HD3	2:B:7489:HOH:O	2.07	0.55
1:F:12:GLU:O	1:F:16:LEU:HD22	2.07	0.55
1:F:177:THR:HB	1:F:203:LYS:HD2	1.88	0.54
1:C:86:HIS:C	2:C:7513:HOH:O	2.46	0.54
1:B:238:ARG:HB3	1:B:238:ARG:CZ	2.36	0.54
1:D:154:HIS:O	1:D:155:ALA:HB3	2.07	0.54
1:F:225:PRO:HG3	2:F:7127:HOH:O	2.07	0.54
1:D:228:GLU:CG	2:D:7338:HOH:O	2.56	0.54
1:D:51:ASN:ND2	2:D:7215:HOH:O	2.41	0.54
1:A:108:GLN:NE2	1:A:108:GLN:HA	2.23	0.53
1:D:197:MSE:HE3	1:D:215:LYS:HG2	1.90	0.53
1:E:190:PRO:CG	1:E:204:ILE:HG22	2.38	0.53
1:E:156:THR:HG23	1:F:137:ASP:OD2	2.08	0.53
1:A:115:TYR:O	1:A:119:GLN:HG2	2.08	0.53
1:D:137:ASP:OD1	1:F:137:ASP:OD1	2.26	0.53
1:E:28:HIS:HA	1:E:32:LEU:HB2	1.89	0.53
1:F:197:MSE:CE	1:F:215:LYS:NZ	2.72	0.53
1:A:143:ALA:HB3	1:A:162:GLU:HG3	1.90	0.52
1:A:15:GLU:HG2	2:A:7506:HOH:O	2.09	0.52
1:C:94:TRP:O	1:C:97:PRO:HD2	2.09	0.52
1:B:238:ARG:CD	2:B:7540:HOH:O	2.56	0.52
1:C:51:ASN:OD1	1:C:53:ILE:N	2.43	0.52
1:B:162:GLU:N	1:B:189:HIS:CE1	2.77	0.52
1:D:89:PRO:CD	2:D:7533:HOH:O	2.54	0.52
1:E:88:ASP:O	1:E:88:ASP:OD1	2.27	0.52
1:F:231:THR:HG22	1:F:240:VAL:CG1	2.39	0.52
1:E:133:SER:O	1:E:137:ASP:HA	2.10	0.52
1:B:238:ARG:CG	2:B:7540:HOH:O	2.57	0.51
1:A:173:LEU:O	1:A:176:VAL:HG13	2.10	0.51
1:C:149:GLY:O	1:C:169:ASP:HA	2.10	0.51
1:E:194:GLU:OE2	1:E:213:TYR:OH	2.19	0.51
1:C:64:GLU:HG2	2:C:7486:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:ASP:OD2	1:F:137:ASP:OD1	2.28	0.51
1:A:88:ASP:O	1:A:90:ALA:N	2.43	0.51
1:F:31:ILE:CD1	1:F:99:LEU:HD21	2.41	0.51
1:D:137:ASP:CG	1:F:137:ASP:OD1	2.49	0.51
1:F:108:GLN:HE22	1:F:111:ARG:HE	1.58	0.51
1:A:8:HIS:O	1:A:12:GLU:HG3	2.11	0.50
1:F:16:LEU:HB3	1:F:27:PHE:CE1	2.46	0.50
1:C:21:PRO:CG	2:C:7555:HOH:O	2.48	0.50
1:B:1:MSE:O	1:B:5:VAL:HG23	2.12	0.50
1:C:87:ARG:NH1	1:C:87:ARG:HG3	2.27	0.50
1:D:109:SER:HB3	1:D:132:ILE:HD11	1.93	0.49
1:D:137:ASP:OD1	1:F:156:THR:HG23	2.12	0.49
1:D:62:ILE:HD13	1:D:124:LEU:HD11	1.95	0.49
1:B:16:LEU:HD21	1:B:93:LEU:HD13	1.94	0.49
1:B:224:ASN:ND2	2:B:7387:HOH:O	2.40	0.49
1:F:109:SER:HB3	1:F:132:ILE:CD1	2.42	0.49
1:A:193:ARG:O	1:A:196:VAL:HG23	2.12	0.49
1:B:173:LEU:O	1:B:176:VAL:HG13	2.13	0.49
1:F:51:ASN:OD1	1:F:53:ILE:N	2.45	0.49
1:A:108:GLN:HE22	1:A:111:ARG:HE	1.60	0.49
1:A:197:MSE:HG3	2:A:7112:HOH:O	2.13	0.48
1:C:177:THR:HB	1:C:203:LYS:HD2	1.94	0.48
1:C:13:ALA:HB2	1:C:31:ILE:HG21	1.94	0.48
1:D:108:GLN:NE2	1:D:108:GLN:HA	2.28	0.48
1:B:197:MSE:HG3	2:B:7414:HOH:O	2.12	0.48
1:B:61:GLU:OE1	2:B:7384:HOH:O	2.20	0.48
1:C:21:PRO:CD	2:C:7555:HOH:O	2.61	0.48
1:A:186:GLY:O	1:A:188:ARG:HG2	2.13	0.48
1:C:62:ILE:HD13	1:C:124:LEU:HD11	1.95	0.48
1:A:63:ILE:HG23	1:A:112:ILE:CD1	2.44	0.48
1:D:197:MSE:CE	1:D:215:LYS:HE2	2.43	0.48
1:A:228:GLU:CG	2:A:7018:HOH:O	2.62	0.47
1:A:57:ILE:HD11	1:D:60:ARG:HD2	1.95	0.47
1:A:89:PRO:CD	2:A:7552:HOH:O	2.57	0.47
1:D:153:ASP:O	1:D:154:HIS:HB2	2.15	0.47
1:D:51:ASN:HB2	1:D:52:PRO:CD	2.44	0.47
1:E:128:LEU:O	1:E:132:ILE:HG13	2.14	0.47
1:E:51:ASN:HB2	1:E:52:PRO:CD	2.45	0.47
1:F:195:GLY:HA3	1:F:213:TYR:CE2	2.49	0.47
1:F:227:PRO:HD2	1:F:239:ILE:HD12	1.95	0.47
1:C:183:LYS:HA	2:C:7568:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:SER:HA	1:E:25:SER:HA	1.96	0.47
1:A:130:ASN:HB3	2:C:7163:HOH:O	2.13	0.47
1:B:190:PRO:HG3	1:B:204:ILE:HG22	1.95	0.47
1:D:190:PRO:C	1:D:191:LYS:HD2	2.35	0.47
1:A:102:LYS:HE3	1:A:136:PHE:O	2.15	0.47
1:A:176:VAL:HA	1:A:202:ALA:O	2.15	0.47
1:D:197:MSE:HE1	1:D:215:LYS:CE	2.44	0.47
1:D:159:VAL:HG11	1:F:154:HIS:CD2	2.50	0.47
1:F:227:PRO:HD2	1:F:239:ILE:CD1	2.45	0.47
1:E:111:ARG:HG2	1:E:111:ARG:HH11	1.80	0.46
1:F:108:GLN:HA	1:F:108:GLN:NE2	2.30	0.46
1:F:96:THR:HB	1:F:97:PRO:HD3	1.96	0.46
1:A:51:ASN:HB2	1:A:52:PRO:CD	2.45	0.46
1:C:95:SER:O	1:C:99:LEU:HG	2.15	0.46
1:E:96:THR:HB	1:E:97:PRO:HD3	1.97	0.46
1:B:92:GLU:CG	1:B:93:LEU:N	2.66	0.46
1:E:82:GLN:HG3	1:E:94:TRP:HZ3	1.80	0.46
1:A:63:ILE:HG23	1:A:112:ILE:HD11	1.97	0.46
1:E:5:VAL:O	1:E:9:ILE:HG13	2.16	0.46
1:F:11:GLN:O	1:F:15:GLU:HG2	2.15	0.46
1:B:176:VAL:HA	1:B:202:ALA:O	2.15	0.46
1:E:32:LEU:HA	1:E:32:LEU:HD23	1.78	0.46
1:E:59:LEU:O	1:E:63:ILE:HG13	2.16	0.46
1:A:150:ILE:HG22	1:A:151:MSE:N	2.31	0.45
1:D:176:VAL:HA	1:D:202:ALA:O	2.16	0.45
1:F:140:ILE:HD11	1:F:152:PHE:HE1	1.81	0.45
1:D:137:ASP:OD1	1:F:156:THR:CG2	2.64	0.45
1:B:115:TYR:O	1:B:119:GLN:HG2	2.16	0.45
1:A:89:PRO:N	2:A:7552:HOH:O	2.49	0.45
1:C:150:ILE:O	2:C:7310:HOH:O	2.21	0.45
1:A:15:GLU:CG	2:A:7400:HOH:O	2.64	0.45
1:B:128:LEU:O	1:B:129:GLN:C	2.54	0.45
1:B:107:ILE:HD11	1:B:152:PHE:CG	2.51	0.45
1:F:92:GLU:HG2	1:F:92:GLU:H	1.55	0.45
1:F:220:SER:OG	1:F:233:ALA:HA	2.17	0.45
1:B:103:GLY:O	1:B:107:ILE:HG13	2.17	0.45
1:B:10:ARG:O	1:B:14:LYS:HG3	2.17	0.45
1:C:28:HIS:HA	1:C:32:LEU:HB2	1.99	0.45
1:C:2:THR:OG1	1:C:75:ASP:OD1	2.31	0.45
1:D:77:ALA:O	1:D:81:ILE:HG13	2.15	0.45
1:E:48:LYS:HE3	1:E:105:HIS:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:117:TRP:CH2	1:F:122:LYS:HD3	2.52	0.44
1:F:87:ARG:HG3	1:F:87:ARG:NH1	2.32	0.44
1:B:239:ILE:HG22	1:B:240:VAL:HG13	1.99	0.44
1:E:173:LEU:O	1:E:176:VAL:HG13	2.17	0.44
1:F:104:PHE:O	1:F:108:GLN:HG2	2.18	0.44
1:F:14:LYS:O	1:F:18:GLU:HG2	2.17	0.44
1:C:29:SER:HB3	1:C:43:TYR:OH	2.17	0.44
1:A:60:ARG:HD2	1:D:57:ILE:HD11	1.99	0.44
1:E:153:ASP:HB3	1:E:173:LEU:HD23	1.98	0.44
1:A:177:THR:HB	1:A:203:LYS:HD2	2.00	0.44
1:C:153:ASP:C	1:C:155:ALA:H	2.21	0.44
1:D:231:THR:HG22	1:D:240:VAL:HG23	1.98	0.44
1:E:63:ILE:HG23	1:E:112:ILE:CD1	2.47	0.44
1:C:44:LEU:O	1:C:48:LYS:HG3	2.17	0.44
1:C:140:ILE:HG23	1:C:160:VAL:HB	2.00	0.44
1:B:92:GLU:HG2	1:B:92:GLU:H	1.61	0.44
1:E:196:VAL:CG1	1:E:197:MSE:N	2.81	0.43
1:A:165:VAL:HG23	1:A:191:LYS:HE3	2.00	0.43
1:A:153:ASP:HB3	1:A:173:LEU:CD2	2.48	0.43
1:F:31:ILE:HG22	1:F:32:LEU:N	2.33	0.43
1:B:108:GLN:CA	1:B:108:GLN:NE2	2.82	0.43
1:C:104:PHE:O	1:C:108:GLN:HG2	2.17	0.43
1:C:31:ILE:HG22	1:C:32:LEU:N	2.32	0.43
1:C:51:ASN:OD1	1:C:51:ASN:C	2.57	0.43
1:D:187:ASP:OD1	1:D:191:LYS:NZ	2.49	0.43
1:E:137:ASP:OD1	1:F:137:ASP:OD1	2.37	0.43
1:E:83:ALA:HB2	1:E:149:GLY:HA2	2.00	0.43
1:F:2:THR:OG1	1:F:75:ASP:OD1	2.29	0.43
1:E:143:ALA:HB3	1:E:162:GLU:HG3	2.00	0.43
1:A:117:TRP:CH2	1:A:122:LYS:HD3	2.54	0.43
1:B:156:THR:CG2	1:C:137:ASP:OD2	2.65	0.43
1:E:20:GLU:HG3	1:E:23:LEU:HB2	2.00	0.43
1:D:90:ALA:N	2:D:7533:HOH:O	2.52	0.43
1:B:51:ASN:HB2	1:B:52:PRO:HD3	2.01	0.43
1:A:53:ILE:CD1	2:A:7567:HOH:O	2.67	0.42
1:C:3:LEU:N	1:C:3:LEU:HD23	2.34	0.42
1:E:152:PHE:CE2	1:E:172:ILE:HD12	2.53	0.42
1:E:182:GLY:O	1:E:184:GLU:N	2.52	0.42
1:E:1:MSE:HB3	1:E:1:MSE:HE3	1.71	0.42
1:F:176:VAL:HA	1:F:202:ALA:O	2.19	0.42
1:A:31:ILE:HD11	1:A:99:LEU:HD21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:GLN:HA	1:B:108:GLN:HE21	1.83	0.42
1:A:211:GLY:O	1:A:212:LYS:C	2.57	0.42
1:F:172:ILE:HD13	1:F:198:ILE:HD12	2.02	0.42
1:B:92:GLU:CG	1:B:93:LEU:HG	2.41	0.42
1:C:212:LYS:HG2	1:C:213:TYR:HD1	1.84	0.42
1:F:33:LYS:NZ	2:F:7140:HOH:O	2.27	0.42
1:E:104:PHE:O	1:E:108:GLN:HG2	2.20	0.41
1:E:65:GLU:CD	1:E:69:SER:HG	2.21	0.41
1:F:23:LEU:HA	1:F:23:LEU:HD12	1.88	0.41
1:D:228:GLU:C	1:D:230:ALA:H	2.24	0.41
1:B:117:TRP:O	1:B:120:ASN:N	2.49	0.41
1:B:51:ASN:HB2	1:B:52:PRO:CD	2.50	0.41
1:D:190:PRO:CG	1:D:204:ILE:HG22	2.50	0.41
1:D:5:VAL:HG21	1:D:78:ALA:HB1	2.02	0.41
1:F:145:LYS:HD2	2:F:7162:HOH:O	2.20	0.41
1:B:173:LEU:HA	1:B:173:LEU:HD23	1.86	0.41
1:C:197:MSE:HE3	2:C:7123:HOH:O	2.20	0.41
1:D:153:ASP:HB3	1:D:173:LEU:HD23	2.03	0.41
1:D:173:LEU:HD22	1:D:174:GLN:OE1	2.20	0.41
1:E:113:THR:HG22	1:E:128:LEU:HB2	2.01	0.41
1:F:54:MSE:HB3	1:F:54:MSE:HE3	1.94	0.41
1:A:196:VAL:HG12	1:A:197:MSE:N	2.36	0.41
1:E:137:ASP:OD1	1:E:156:THR:HG23	2.20	0.41
1:B:234:GLY:HA2	2:B:7497:HOH:O	2.21	0.41
1:D:96:THR:O	1:D:97:PRO:C	2.58	0.41
1:B:228:GLU:HG3	2:B:7398:HOH:O	2.19	0.41
1:C:12:GLU:OE1	1:C:94:TRP:HD1	2.03	0.41
1:D:224:ASN:HA	1:D:225:PRO:HD3	1.95	0.41
1:E:24:ALA:O	1:E:25:SER:C	2.58	0.41
1:B:63:ILE:HG23	1:B:112:ILE:CD1	2.51	0.41
1:D:194:GLU:O	1:D:212:LYS:HA	2.21	0.41
1:D:1:MSE:CB	2:D:7468:HOH:O	2.69	0.41
1:C:141:HIS:HB3	1:C:162:GLU:HA	2.02	0.41
1:C:179:GLY:O	1:C:190:PRO:HD3	2.21	0.41
1:F:149:GLY:O	1:F:169:ASP:HA	2.20	0.41
1:E:226:VAL:HA	1:E:227:PRO:HD3	1.81	0.40
1:E:46:ALA:O	1:E:50:ALA:HB2	2.21	0.40
1:D:97:PRO:HA	1:D:101:LEU:HG	2.02	0.40
1:F:63:ILE:HG23	1:F:112:ILE:CD1	2.52	0.40
1:A:145:LYS:HG3	2:A:7345:HOH:O	2.21	0.40
1:C:32:LEU:HA	1:C:32:LEU:HD23	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:153:ASP:O	1:E:155:ALA:N	2.48	0.40
1:C:184:GLU:N	2:C:7568:HOH:O	2.38	0.40
1:D:191:LYS:N	1:D:191:LYS:HD2	2.36	0.40
1:F:69:SER:HB2	1:F:115:TYR:CE1	2.56	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	238/242 (98%)	229 (96%)	8 (3%)	1 (0%)	34	29
1	B	238/242 (98%)	225 (94%)	13 (6%)	0	100	100
1	C	238/242 (98%)	229 (96%)	9 (4%)	0	100	100
1	D	238/242 (98%)	224 (94%)	13 (6%)	1 (0%)	34	29
1	E	238/242 (98%)	225 (94%)	11 (5%)	2 (1%)	19	12
1	F	238/242 (98%)	227 (95%)	11 (5%)	0	100	100
All	All	1428/1452 (98%)	1359 (95%)	65 (5%)	4 (0%)	41	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	229	TYR
1	E	183	LYS
1	E	89	PRO
1	A	89	PRO

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	190/192 (99%)	183 (96%)	7 (4%)	34	32
1	B	191/192 (100%)	184 (96%)	7 (4%)	34	32
1	C	189/192 (98%)	181 (96%)	8 (4%)	30	28
1	D	190/192 (99%)	183 (96%)	7 (4%)	34	32
1	E	192/192 (100%)	183 (95%)	9 (5%)	26	23
1	F	190/192 (99%)	179 (94%)	11 (6%)	20	15
All	All	1142/1152 (99%)	1093 (96%)	49 (4%)	29	27

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	21	PRO
1	A	110	TYR
1	A	137	ASP
1	A	176	VAL
1	A	188	ARG
1	A	227	PRO
1	B	16	LEU
1	B	18	GLU
1	B	35	GLN
1	B	82	GLN
1	B	110	TYR
1	B	120	ASN
1	B	238	ARG
1	C	16	LEU
1	C	72	SER
1	C	110	TYR
1	C	137	ASP
1	C	176	VAL
1	C	191	LYS
1	C	194	GLU
1	C	238	ARG

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Mol	Chain	Res	Type
1	D	16	LEU
1	D	110	TYR
1	D	115	TYR
1	D	137	ASP
1	D	176	VAL
1	D	188	ARG
1	D	212	LYS
1	E	1	MSE
1	E	16	LEU
1	E	35	GLN
1	E	68	GLN
1	E	86	HIS
1	E	110	TYR
1	E	176	VAL
1	E	194	GLU
1	E	238	ARG
1	F	11	GLN
1	F	16	LEU
1	F	58	SER
1	F	68	GLN
1	F	82	GLN
1	F	92	GLU
1	F	110	TYR
1	F	115	TYR
1	F	176	VAL
1	F	187	ASP
1	F	194	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	86	HIS
1	A	108	GLN
1	A	141	HIS
1	B	7	GLN
1	B	108	GLN
1	B	141	HIS
1	B	224	ASN
1	C	7	GLN
1	C	19	ASN
1	C	108	GLN

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Mol	Chain	Res	Type
1	C	141	HIS
1	C	148	HIS
1	C	154	HIS
1	D	7	GLN
1	D	68	GLN
1	D	141	HIS
1	E	19	ASN
1	E	35	GLN
1	E	108	GLN
1	E	141	HIS
1	F	7	GLN
1	F	82	GLN
1	F	108	GLN
1	F	141	HIS
1	F	154	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 carbohydrates 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	235/242 (97%)	0.04	11 (4%) 31 41	16, 25, 41, 56	0
1	B	235/242 (97%)	0.05	7 (2%) 50 59	16, 25, 41, 55	0
1	C	235/242 (97%)	0.09	7 (2%) 50 59	17, 25, 42, 52	0
1	D	235/242 (97%)	0.05	8 (3%) 45 53	15, 24, 40, 50	0
1	E	235/242 (97%)	0.12	11 (4%) 31 41	16, 25, 42, 52	0
1	F	235/242 (97%)	-0.02	9 (3%) 40 49	16, 23, 41, 53	0
All	All	1410/1452 (97%)	0.05	53 (3%) 40 49	15, 25, 42, 56	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	184	GLU	5.8
1	B	185	SER	5.7
1	C	185	SER	5.6
1	A	184	GLU	5.5
1	E	184	GLU	5.5
1	C	182	GLY	5.4
1	E	185	SER	5.3
1	D	186	GLY	5.1
1	C	184	GLU	5.0
1	D	89	PRO	4.9
1	C	183	LYS	4.7
1	B	89	PRO	4.5
1	C	186	GLY	4.4
1	A	89	PRO	4.2
1	E	182	GLY	3.9
1	F	90	ALA	3.9
1	A	183	LYS	3.8
1	F	184	GLU	3.7
1	A	86	HIS	3.7

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Mol	Chain	Res	Type	RSRZ
1	F	185	SER	3.4
1	E	89	PRO	3.3
1	F	89	PRO	3.1
1	F	182	GLY	3.1
1	A	90	ALA	3.1
1	F	183	LYS	3.1
1	E	86	HIS	3.0
1	A	186	GLY	3.0
1	B	182	GLY	3.0
1	E	186	GLY	3.0
1	E	4	ASP	3.0
1	A	238	ARG	2.8
1	C	94	TRP	2.7
1	E	183	LYS	2.5
1	D	182	GLY	2.5
1	F	85	ARG	2.4
1	A	240	VAL	2.4
1	D	184	GLU	2.4
1	D	213	TYR	2.3
1	C	89	PRO	2.3
1	F	92	GLU	2.3
1	B	229	TYR	2.3
1	A	92	GLU	2.3
1	F	86	HIS	2.3
1	E	90	ALA	2.3
1	D	238	ARG	2.3
1	B	240	VAL	2.3
1	E	3	LEU	2.2
1	A	229	TYR	2.2
1	A	182	GLY	2.2
1	E	229	TYR	2.1
1	D	3	LEU	2.1
1	D	92	GLU	2.0
1	B	186	GLY	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

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.