Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:40 PM GMT

PDB ID	:	1LGY
Title	:	LIPASE II FROM RHIZOPUS NIVEUS
Authors	:	Kohno, M.; Funatsu, J.; Mikami, B.; Kugimiya, W.; Matsuo, T.; Morita, Y.
Deposited on	:	1996-05-23
Resolution	:	2.20 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 ⓘ symbol.

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

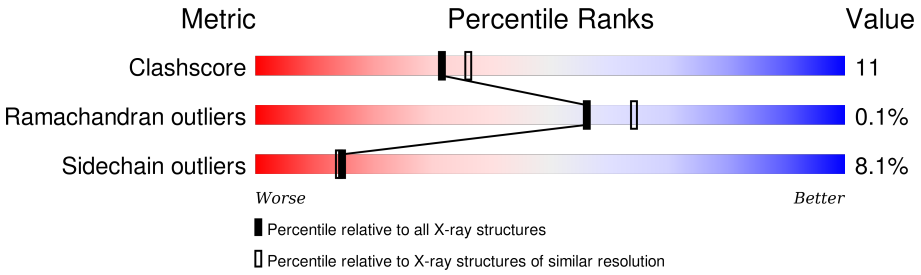
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 2.20 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))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	269	71% 23% 5%
1	B	269	74% 20%
1	C	269	67% 28%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6682 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 TRIACYLGLYCEROL LIPASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2068	1326	347	388	7			
1	B	265	Total	C	N	O	S	0	0	0
			2068	1326	347	388	7			
1	C	265	Total	C	N	O	S	0	0	0
			2068	1326	347	388	7			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	135	Total	O	0	0
			135	135		
2	B	157	Total	O	0	0
			157	157		
2	C	186	Total	O	0	0
			186	186		

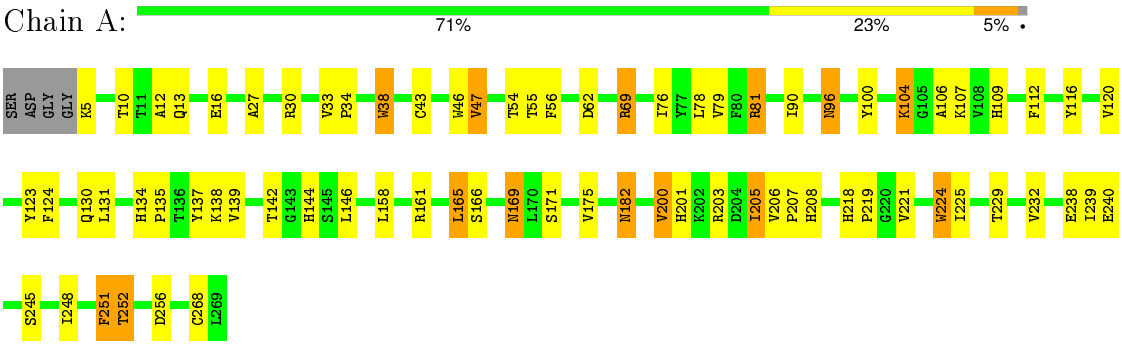
3 Residue-property plots (i)

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.

Note EDS was not executed.

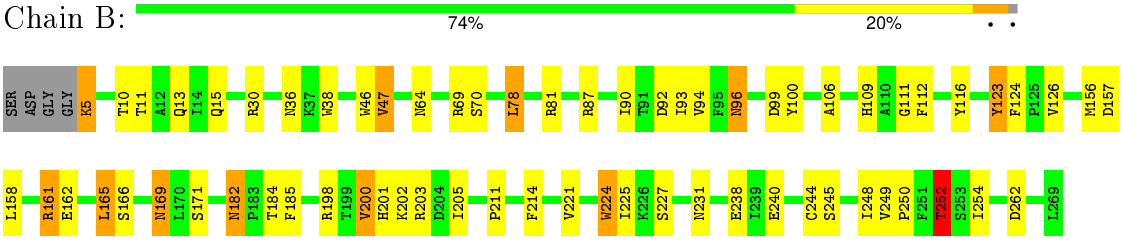
- Molecule 1: TRIACYLGLYCEROL LIPASE

Chain A:



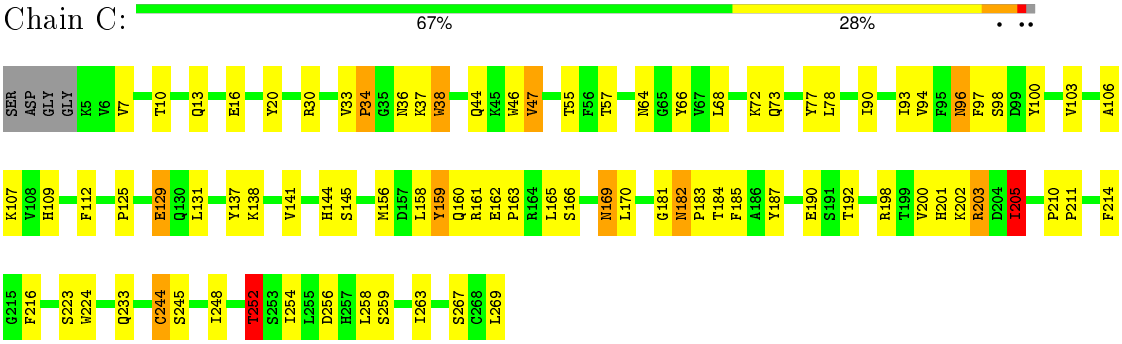
- Molecule 1: TRIACYLGLYCEROL LIPASE

Chain B:



- Molecule 1: TRIACYLGLYCEROL LIPASE

Chain C:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α, β, γ	83.70Å 83.70Å 137.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.186 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6682	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

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	0.91	0/2122	1.56	24/2889 (0.8%)
1	B	0.87	0/2122	1.57	24/2889 (0.8%)
1	C	0.82	0/2122	1.53	26/2889 (0.9%)
All	All	0.87	0/6366	1.55	74/8667 (0.9%)

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

There are no bond length outliers.

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	161	ARG	NE-CZ-NH2	-13.93	113.33	120.30
1	A	38	TRP	CD1-CG-CD2	8.86	113.39	106.30
1	C	205	ILE	CB-CA-C	-8.74	94.12	111.60
1	B	38	TRP	CD1-CG-CD2	8.62	113.20	106.30
1	C	38	TRP	CD1-CG-CD2	8.51	113.11	106.30
1	C	38	TRP	CE2-CD2-CG	-8.20	100.74	107.30
1	A	104	LYS	CA-CB-CG	8.10	131.22	113.40
1	B	38	TRP	CE2-CD2-CG	-8.05	100.86	107.30
1	C	224	TRP	CD1-CG-CD2	7.98	112.69	106.30
1	B	38	TRP	CG-CD2-CE3	7.85	140.96	133.90
1	B	46	TRP	CD1-CG-CD2	7.81	112.55	106.30
1	A	38	TRP	CE2-CD2-CG	-7.61	101.21	107.30
1	B	100	TYR	CB-CG-CD2	-7.52	116.49	121.00
1	C	46	TRP	CE2-CD2-CG	-7.50	101.30	107.30
1	A	161	ARG	NE-CZ-NH1	7.45	124.03	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	46	TRP	CD1-CG-CD2	7.39	112.22	106.30
1	A	205	ILE	CB-CA-C	-7.35	96.90	111.60
1	B	161	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	B	38	TRP	CB-CG-CD1	-7.33	117.47	127.00
1	C	224	TRP	CE2-CD2-CG	-7.30	101.46	107.30
1	B	224	TRP	CE2-CD2-CG	-7.22	101.52	107.30
1	A	46	TRP	CD1-CG-CD2	7.20	112.06	106.30
1	B	156	MET	CG-SD-CE	7.19	111.71	100.20
1	A	46	TRP	CE2-CD2-CG	-7.00	101.70	107.30
1	B	46	TRP	CE2-CD2-CG	-6.93	101.75	107.30
1	C	198	ARG	NE-CZ-NH1	6.91	123.76	120.30
1	B	69	ARG	NE-CZ-NH1	6.87	123.74	120.30
1	B	224	TRP	CG-CD2-CE3	6.85	140.07	133.90
1	A	137	TYR	CB-CG-CD2	-6.67	117.00	121.00
1	A	224	TRP	CD1-CG-CD2	6.51	111.51	106.30
1	B	224	TRP	CD1-CG-CD2	6.50	111.50	106.30
1	C	46	TRP	CG-CD2-CE3	6.46	139.72	133.90
1	C	252	THR	N-CA-CB	-6.36	98.21	110.30
1	A	224	TRP	CE2-CD2-CG	-6.34	102.23	107.30
1	A	116	TYR	CB-CG-CD2	-6.24	117.26	121.00
1	A	30	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	C	141	VAL	CG1-CB-CG2	-6.17	101.03	110.90
1	B	30	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	B	30	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	C	161	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	C	159	TYR	CB-CG-CD2	-5.96	117.42	121.00
1	C	203	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	C	46	TRP	CB-CG-CD1	-5.85	119.39	127.00
1	B	252	THR	N-CA-CB	-5.80	99.28	110.30
1	B	240	GLU	N-CA-CB	-5.78	100.19	110.60
1	A	120	VAL	CG1-CB-CG2	-5.68	101.81	110.90
1	A	38	TRP	CG-CD1-NE1	-5.65	104.45	110.10
1	C	161	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	C	244	CYS	CA-C-N	5.56	129.43	117.20
1	B	224	TRP	CB-CG-CD1	-5.56	119.77	127.00
1	B	171	SER	N-CA-CB	-5.54	102.19	110.50
1	A	218	HIS	CA-CB-CG	5.54	123.02	113.60
1	C	205	ILE	CB-CG1-CD1	-5.53	98.41	113.90
1	C	30	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	A	229	THR	N-CA-CB	-5.52	99.81	110.30
1	C	100	TYR	CB-CG-CD2	-5.52	117.69	121.00
1	B	203	ARG	NE-CZ-NH2	-5.50	117.55	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	43	CYS	CA-C-N	-5.47	105.16	117.20
1	A	62	ASP	CB-CG-OD1	5.43	123.19	118.30
1	C	205	ILE	N-CA-CB	5.41	123.24	110.80
1	C	30	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	A	81	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	A	38	TRP	CG-CD2-CE3	5.30	138.67	133.90
1	B	126	VAL	CG1-CB-CG2	-5.23	102.54	110.90
1	C	224	TRP	CG-CD2-CE3	5.23	138.60	133.90
1	A	46	TRP	CG-CD2-CE3	5.18	138.56	133.90
1	B	38	TRP	CG-CD1-NE1	-5.17	104.94	110.10
1	A	240	GLU	CA-C-N	-5.16	105.84	117.20
1	C	203	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	69	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	B	244	CYS	CA-C-N	5.06	128.34	117.20
1	C	38	TRP	CG-CD2-CE3	5.04	138.44	133.90
1	B	203	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	C	137	TYR	CB-CG-CD2	-5.02	117.99	121.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	116	TYR	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	2068	0	2035	42	0
1	B	2068	0	2035	37	0
1	C	2068	0	2035	53	0
2	A	135	0	0	0	0
2	B	157	0	0	4	0
2	C	186	0	0	3	0
All	All	6682	0	6105	132	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 (132) 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:C:267:SER:HB3	1:C:269:LEU:HG	1.61	0.83
1:C:97:PHE:HD2	1:C:107:LYS:HB3	1.54	0.73
1:C:205:ILE:HG23	1:C:252:THR:HG23	1.72	0.71
1:A:201:HIS:HB2	1:A:225:ILE:HD12	1.73	0.71
1:C:38:TRP:HZ2	1:C:47:VAL:HG13	1.58	0.67
1:A:10:THR:HG22	1:A:13:GLN:HG3	1.77	0.67
1:A:96:ASN:H	1:A:96:ASN:HD22	1.44	0.65
1:C:97:PHE:CD2	1:C:107:LYS:HB3	2.30	0.65
1:A:203:ARG:NH1	1:A:248:ILE:HB	2.12	0.64
1:A:38:TRP:HZ2	1:A:47:VAL:HG13	1.63	0.63
1:C:156:MET:SD	1:C:192:THR:HG21	2.39	0.63
1:B:10:THR:H	1:B:13:GLN:HE21	1.45	0.62
1:C:192:THR:HG22	2:C:600:HOH:O	1.99	0.61
1:C:107:LYS:H	1:C:182:ASN:ND2	1.99	0.60
1:A:169:ASN:HD22	1:A:169:ASN:H	1.49	0.59
1:B:205:ILE:HG23	1:B:252:THR:HG23	1.84	0.59
1:A:109:HIS:HD2	1:A:112:PHE:H	1.50	0.59
1:A:238:GLU:O	1:A:239:ILE:HD13	2.04	0.58
1:B:166:SER:H	1:B:169:ASN:ND2	2.02	0.58
1:B:205:ILE:HG23	1:B:252:THR:CG2	2.33	0.57
1:A:10:THR:H	1:A:13:GLN:HE21	1.53	0.56
1:C:205:ILE:HD11	1:C:254:ILE:HD13	1.87	0.56
1:C:57:THR:HG23	1:C:64:ASN:OD1	2.05	0.56
1:A:79:VAL:HG22	1:A:142:THR:HG22	1.86	0.56
1:B:96:ASN:H	1:B:96:ASN:HD22	1.53	0.54
1:C:109:HIS:HD2	1:C:112:PHE:H	1.54	0.54
1:A:165:LEU:HA	1:A:169:ASN:HD21	1.72	0.54
1:C:55:THR:HG22	1:C:66:TYR:HB3	1.90	0.54
1:B:87:ARG:HD2	2:B:645:HOH:O	2.08	0.54
1:C:10:THR:H	1:C:13:GLN:NE2	2.07	0.53
1:C:267:SER:CB	1:C:269:LEU:HG	2.36	0.53
1:C:68:LEU:HD23	1:C:77:TYR:CD1	2.44	0.53
1:A:47:VAL:HG22	1:A:47:VAL:O	2.10	0.52
1:A:200:VAL:HB	1:A:207:PRO:HG2	1.91	0.52
1:B:78:LEU:HD11	1:B:123:TYR:CD1	2.45	0.52
1:C:106:ALA:HA	1:C:184:THR:HB	1.92	0.52
1:C:125:PRO:O	1:C:129:GLU:HB2	2.10	0.52
1:C:210:PRO:HB2	1:C:216:PHE:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:VAL:HG23	1:A:55:THR:HB	1.92	0.51
1:A:107:LYS:H	1:A:182:ASN:ND2	2.07	0.51
1:B:11:THR:HG22	2:B:750:HOH:O	2.10	0.51
1:C:182:ASN:ND2	1:C:185:PHE:H	2.08	0.51
1:A:10:THR:HG23	1:A:12:ALA:HB3	1.92	0.51
1:C:107:LYS:H	1:C:182:ASN:HD21	1.59	0.51
1:A:205:ILE:HG23	1:A:252:THR:CG2	2.41	0.51
1:C:47:VAL:HB	1:C:73:GLN:HE22	1.76	0.50
1:A:38:TRP:CZ2	1:A:47:VAL:HG13	2.46	0.50
1:C:10:THR:HG22	1:C:13:GLN:NE2	2.26	0.50
1:C:166:SER:H	1:C:169:ASN:ND2	2.09	0.50
1:B:5:LYS:HD3	2:B:474:HOH:O	2.11	0.50
1:A:124:PHE:CE2	1:A:158:LEU:HD23	2.47	0.50
1:B:227:SER:O	1:B:231:ASN:HB3	2.13	0.49
1:B:10:THR:N	1:B:13:GLN:HE21	2.10	0.49
1:C:165:LEU:HA	1:C:169:ASN:HD21	1.77	0.49
1:C:201:HIS:HD2	1:C:256:ASP:O	1.96	0.49
1:B:211:PRO:HG2	1:B:214:PHE:HD2	1.78	0.49
1:C:211:PRO:HG2	1:C:214:PHE:HD2	1.77	0.49
1:C:10:THR:N	1:C:13:GLN:NE2	2.60	0.48
1:A:224:TRP:O	1:A:232:VAL:HA	2.14	0.48
1:A:10:THR:H	1:A:13:GLN:NE2	2.11	0.47
1:A:69:ARG:HD3	1:A:130:GLN:OE1	2.14	0.47
1:B:47:VAL:HG21	1:B:70:SER:HB2	1.95	0.47
1:B:205:ILE:HG12	1:B:252:THR:HG23	1.95	0.47
1:B:78:LEU:HD11	1:B:123:TYR:CE1	2.50	0.47
1:A:10:THR:HG22	1:A:13:GLN:CG	2.45	0.47
1:A:27:ALA:O	1:A:81:ARG:HD3	2.15	0.47
1:C:16:GLU:HG2	1:C:20:TYR:HE2	1.80	0.46
1:B:106:ALA:HA	1:B:184:THR:HB	1.98	0.46
1:C:205:ILE:HD11	1:C:254:ILE:CD1	2.45	0.46
1:C:159:TYR:HE2	2:C:600:HOH:O	1.98	0.46
1:C:169:ASN:HD22	1:C:169:ASN:H	1.64	0.46
1:A:245:SER:O	1:A:248:ILE:HG13	2.15	0.46
1:C:103:VAL:HB	1:C:106:ALA:HB3	1.97	0.46
1:B:166:SER:H	1:B:169:ASN:HD21	1.63	0.46
1:B:124:PHE:CE1	1:B:158:LEU:HD23	2.50	0.46
1:B:165:LEU:HA	1:B:169:ASN:HD21	1.80	0.45
1:B:11:THR:O	1:B:15:GLN:HG3	2.16	0.45
1:A:201:HIS:HD2	1:A:256:ASP:O	1.99	0.45
1:A:56:PHE:CE2	1:A:123:TYR:HB3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ILE:O	1:A:208:HIS:HB2	2.16	0.45
1:C:131:LEU:HD11	1:C:165:LEU:CD2	2.47	0.45
1:B:109:HIS:HD2	1:B:112:PHE:H	1.64	0.45
1:C:38:TRP:CZ2	1:C:47:VAL:HG13	2.45	0.45
1:C:181:GLY:O	1:C:216:PHE:HA	2.16	0.44
1:A:268:CYS:O	1:A:268:CYS:SG	2.74	0.44
1:A:203:ARG:HH11	1:A:203:ARG:HG2	1.83	0.44
1:C:96:ASN:HB3	2:C:768:HOH:O	2.17	0.44
1:C:245:SER:O	1:C:248:ILE:HG13	2.18	0.44
1:B:201:HIS:HB2	1:B:225:ILE:HD12	1.99	0.44
1:B:182:ASN:ND2	1:B:185:PHE:H	2.16	0.44
1:C:159:TYR:HB2	1:C:170:LEU:HD23	2.00	0.44
1:A:76:ILE:O	1:A:139:VAL:HA	2.18	0.44
1:A:169:ASN:HD22	1:A:169:ASN:N	2.11	0.43
1:B:109:HIS:CD2	1:B:111:GLY:H	2.35	0.43
1:B:94:VAL:O	1:B:109:HIS:HE1	2.02	0.43
1:C:259:SER:HA	1:C:263:ILE:O	2.18	0.43
1:C:109:HIS:HD2	1:C:112:PHE:N	2.16	0.43
1:B:202:LYS:HG2	1:B:225:ILE:O	2.19	0.42
1:B:64:ASN:HB2	1:B:81:ARG:HG3	2.02	0.42
1:A:5:LYS:HA	1:A:5:LYS:HD2	1.75	0.42
1:A:205:ILE:HG23	1:A:252:THR:HG21	2.00	0.42
1:B:198:ARG:HD2	2:B:439:HOH:O	2.19	0.42
1:B:94:VAL:O	1:B:109:HIS:CE1	2.73	0.42
1:C:138:LYS:HB3	1:C:138:LYS:HE3	1.77	0.42
1:A:203:ARG:NH1	1:A:251:PHE:HB2	2.35	0.42
1:A:166:SER:H	1:A:169:ASN:ND2	2.18	0.41
1:B:200:VAL:O	1:B:224:TRP:HD1	2.03	0.41
1:B:93:ILE:HG21	1:B:93:ILE:HD13	1.83	0.41
1:C:90:ILE:O	1:C:93:ILE:HG22	2.19	0.41
1:C:203:ARG:HB3	1:C:252:THR:HA	2.02	0.41
1:B:90:ILE:HD13	1:B:90:ILE:HA	1.90	0.41
1:C:47:VAL:HG22	1:C:47:VAL:O	2.21	0.41
1:B:245:SER:O	1:B:248:ILE:HG13	2.20	0.41
1:A:134:HIS:HA	1:A:135:PRO:HD3	1.85	0.41
1:B:157:ASP:OD1	1:B:161:ARG:HD3	2.21	0.41
1:C:158:LEU:O	1:C:162:GLU:HB3	2.21	0.41
1:C:131:LEU:HD11	1:C:165:LEU:HD21	2.03	0.41
1:C:233:GLN:NE2	1:C:244:CYS:SG	2.91	0.41
1:B:166:SER:N	1:B:169:ASN:HD21	2.19	0.41
1:A:144:HIS:HA	1:A:175:VAL:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:ILE:CD1	1:A:131:LEU:HD23	2.51	0.41
1:C:144:HIS:CE1	1:C:145:SER:HB2	2.55	0.41
1:C:183:PRO:O	1:C:187:TYR:CD2	2.74	0.41
1:C:33:VAL:HA	1:C:34:PRO:HA	1.89	0.41
1:C:10:THR:HG23	1:C:13:GLN:H	1.86	0.40
1:A:203:ARG:HH12	1:A:248:ILE:HB	1.84	0.40
1:B:205:ILE:HG21	1:B:205:ILE:HD13	1.86	0.40
1:A:90:ILE:HD11	1:A:206:VAL:HG22	2.03	0.40
1:C:156:MET:O	1:C:160:GLN:HG3	2.21	0.40
1:C:94:VAL:O	1:C:109:HIS:CE1	2.74	0.40
1:B:249:VAL:HA	1:B:250:PRO:HA	1.80	0.40
1:A:100:TYR:HB3	1:A:106:ALA:HB3	2.04	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	263/269 (98%)	245 (93%)	18 (7%)	0	100	100
1	B	263/269 (98%)	250 (95%)	13 (5%)	0	100	100
1	C	263/269 (98%)	242 (92%)	20 (8%)	1 (0%)	39	42
All	All	789/807 (98%)	737 (93%)	51 (6%)	1 (0%)	56	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	202	LYS

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	231/233 (99%)	213 (92%)	18 (8%)	16	15
1	B	231/233 (99%)	213 (92%)	18 (8%)	16	15
1	C	231/233 (99%)	211 (91%)	20 (9%)	13	12
All	All	693/699 (99%)	637 (92%)	56 (8%)	15	14

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	34	PRO
1	A	47	VAL
1	A	54	THR
1	A	78	LEU
1	A	96	ASN
1	A	104	LYS
1	A	138	LYS
1	A	146	LEU
1	A	165	LEU
1	A	169	ASN
1	A	171	SER
1	A	182	ASN
1	A	200	VAL
1	A	219	PRO
1	A	221	VAL
1	A	251	PHE
1	A	252	THR
1	B	5	LYS
1	B	36	ASN
1	B	47	VAL
1	B	78	LEU
1	B	92	ASP
1	B	96	ASN
1	B	99	ASP
1	B	123	TYR

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Mol	Chain	Res	Type
1	B	162	GLU
1	B	165	LEU
1	B	169	ASN
1	B	182	ASN
1	B	200	VAL
1	B	221	VAL
1	B	238	GLU
1	B	252	THR
1	B	254	ILE
1	B	262	ASP
1	C	7	VAL
1	C	34	PRO
1	C	36	ASN
1	C	37	LYS
1	C	44	GLN
1	C	47	VAL
1	C	72	LYS
1	C	78	LEU
1	C	96	ASN
1	C	98	SER
1	C	129	GLU
1	C	163	PRO
1	C	169	ASN
1	C	182	ASN
1	C	190	GLU
1	C	200	VAL
1	C	205	ILE
1	C	223	SER
1	C	252	THR
1	C	258	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	96	ASN
1	A	109	HIS
1	A	118	GLN
1	A	169	ASN
1	A	182	ASN
1	A	201	HIS
1	A	212	GLN

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Mol	Chain	Res	Type
1	B	13	GLN
1	B	73	GLN
1	B	96	ASN
1	B	109	HIS
1	B	118	GLN
1	B	169	ASN
1	B	182	ASN
1	B	212	GLN
1	C	13	GLN
1	C	36	ASN
1	C	44	GLN
1	C	73	GLN
1	C	96	ASN
1	C	109	HIS
1	C	169	ASN
1	C	182	ASN
1	C	201	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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.