Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 10:22 PM GMT

PDB ID	:	1TAH
Title	:	THE CRYSTAL STRUCTURE OF TRIACYLGLYCEROL LIPASE FROM PSEUDOMONAS GLUMAE REVEALS A PARTIALLY REDUNDANT CATALYTIC ASPARTATE
Authors	:	Noble, M.E.M.; Cleasby, A.; Johnson, L.N.; Egmond, M.; Frenken, L.G.J.
Deposited on	:	1993-12-21
Resolution	:	3.00 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:

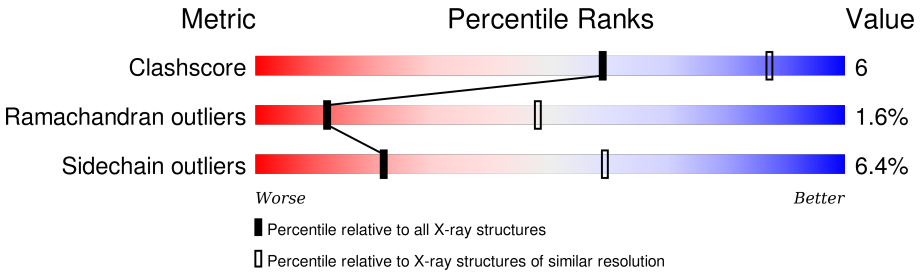
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 3.00 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	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)

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	318	83% 14%
1	B	318	81% 16%
1	C	318	84% 12%
1	D	318	80% 17%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9320 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 LIPASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	318	Total	C	N	O	S	0	0	0
			2329	1450	410	466	3			
1	A	318	Total	C	N	O	S	0	0	0
			2329	1450	410	466	3			
1	C	318	Total	C	N	O	S	0	0	0
			2329	1450	410	466	3			
1	D	318	Total	C	N	O	S	0	0	0
			2329	1450	410	466	3			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		
2	D	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		

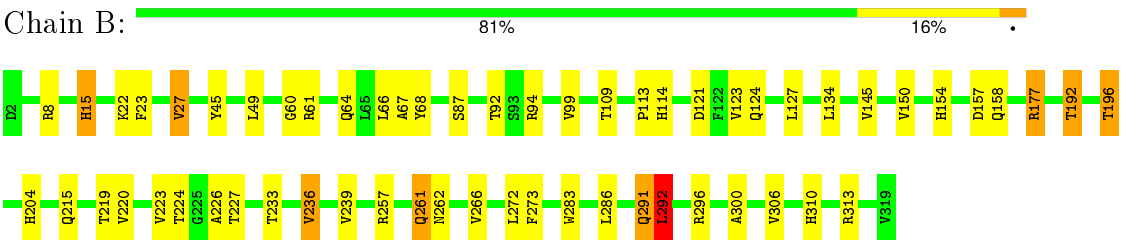
3 Residue-property plots (i)

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.

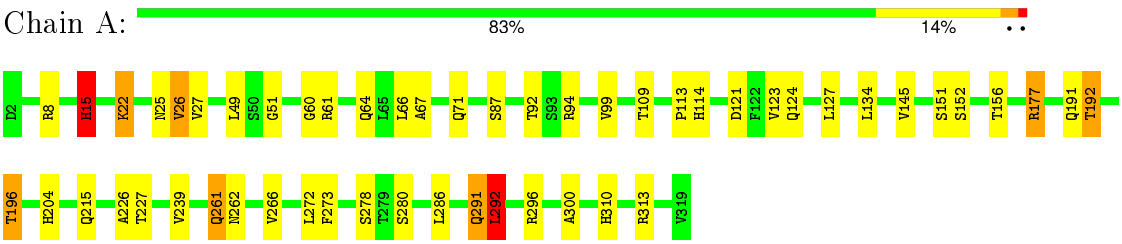
- Molecule 1: LIPASE

Chain B:



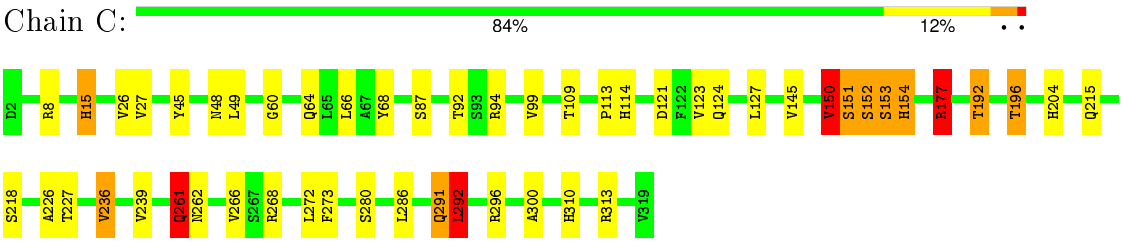
- Molecule 1: LIPASE

Chain A:



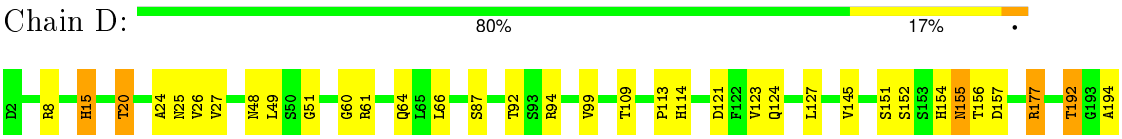
- Molecule 1: LIPASE

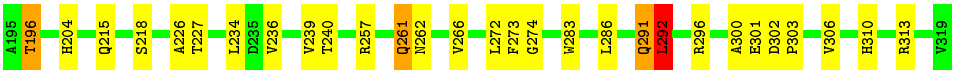
Chain C:



- Molecule 1: LIPASE

Chain D:




A195
T196
H204
Q215
S218
A226
T227
L234
D235
V236
V239
T240
R257
Q261
N262
V266
L272
F273
G274
W283
L286
Q291
L292
R296
A300
E301
D302
P303
V306
H310
R313
V319

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 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	158.16Å 158.64Å 63.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.159 , 0.226	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9320	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0/2371	1.24	15/3242 (0.5%)
1	B	0.58	0/2371	1.29	16/3242 (0.5%)
1	C	0.59	0/2371	1.31	22/3242 (0.7%)
1	D	0.58	0/2371	1.26	13/3242 (0.4%)
All	All	0.58	0/9484	1.28	66/12968 (0.5%)

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	C	0	1

There are no bond length outliers.

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	291	GLN	CA-C-N	-9.74	95.78	117.20
1	D	291	GLN	CA-C-N	-9.72	95.81	117.20
1	C	291	GLN	CA-C-N	-9.65	95.98	117.20
1	A	291	GLN	CA-C-N	-9.46	96.40	117.20
1	A	291	GLN	O-C-N	8.79	136.76	122.70
1	C	291	GLN	O-C-N	8.78	136.75	122.70
1	B	291	GLN	O-C-N	8.56	136.40	122.70
1	C	150	VAL	CA-CB-CG2	-8.52	98.12	110.90
1	D	291	GLN	O-C-N	8.34	136.04	122.70
1	C	236	VAL	CA-CB-CG1	-8.15	98.68	110.90
1	C	152	SER	CA-C-N	-7.63	100.42	117.20
1	B	177	ARG	NE-CZ-NH1	7.24	123.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	292	LEU	CA-C-N	-7.19	101.39	117.20
1	C	150	VAL	N-CA-CB	-7.11	95.86	111.50
1	C	292	LEU	CA-C-N	-7.11	101.56	117.20
1	B	292	LEU	CA-C-N	-7.08	101.63	117.20
1	A	292	LEU	CA-C-N	-7.07	101.65	117.20
1	A	177	ARG	NE-CZ-NH1	6.80	123.70	120.30
1	D	177	ARG	NE-CZ-NH1	6.78	123.69	120.30
1	C	291	GLN	C-N-CA	6.61	138.23	121.70
1	A	291	GLN	C-N-CA	6.55	138.07	121.70
1	B	291	GLN	C-N-CA	6.53	138.02	121.70
1	C	153	SER	N-CA-C	-6.43	93.63	111.00
1	D	291	GLN	C-N-CA	6.42	137.74	121.70
1	D	61	ARG	NE-CZ-NH2	-6.41	117.09	120.30
1	B	27	VAL	CA-CB-CG1	-6.07	101.79	110.90
1	C	177	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	C	150	VAL	CA-CB-CG1	6.04	119.96	110.90
1	D	8	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	B	220	VAL	CA-C-N	-5.90	104.22	117.20
1	C	268	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	A	99	VAL	CA-CB-CG2	-5.82	102.17	110.90
1	B	261	GLN	N-CA-CB	5.80	121.03	110.60
1	A	61	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	D	261	GLN	N-CA-CB	5.77	120.99	110.60
1	C	150	VAL	O-C-N	-5.71	113.56	122.70
1	C	145	VAL	CA-CB-CG2	-5.71	102.34	110.90
1	C	99	VAL	CA-CB-CG2	-5.67	102.39	110.90
1	A	261	GLN	N-CA-CB	5.64	120.74	110.60
1	C	261	GLN	N-CA-CB	5.63	120.73	110.60
1	B	261	GLN	N-CA-C	-5.57	95.96	111.00
1	B	145	VAL	CA-CB-CG2	-5.56	102.56	110.90
1	B	61	ARG	NE-CZ-NH2	-5.55	117.53	120.30
1	C	152	SER	O-C-N	5.54	131.57	122.70
1	B	8	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	8	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	D	145	VAL	CA-CB-CG2	-5.52	102.62	110.90
1	C	261	GLN	N-CA-C	-5.52	96.09	111.00
1	B	257	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	D	99	VAL	CA-CB-CG2	-5.49	102.66	110.90
1	C	236	VAL	CA-CB-CG2	5.49	119.13	110.90
1	A	261	GLN	N-CA-C	-5.43	96.33	111.00
1	B	99	VAL	CA-CB-CG2	-5.43	102.76	110.90
1	C	8	ARG	NE-CZ-NH1	5.41	123.00	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	150	VAL	N-CA-C	5.41	125.59	111.00
1	A	145	VAL	CA-CB-CG2	-5.40	102.80	110.90
1	D	261	GLN	N-CA-C	-5.27	96.77	111.00
1	A	152	SER	CA-C-N	-5.25	105.65	117.20
1	D	257	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	A	26	VAL	CA-CB-CG2	-5.19	103.12	110.90
1	A	292	LEU	N-CA-C	-5.19	97.00	111.00
1	A	15	HIS	CA-CB-CG	5.15	122.36	113.60
1	C	292	LEU	N-CA-C	-5.15	97.10	111.00
1	D	261	GLN	CB-CA-C	-5.11	100.18	110.40
1	B	236	VAL	CA-CB-CG2	-5.07	103.30	110.90
1	B	292	LEU	N-CA-C	-5.01	97.48	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	154	HIS	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	2329	0	2282	25	0
1	B	2329	0	2282	27	0
1	C	2329	0	2282	30	0
1	D	2329	0	2282	29	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
All	All	9320	0	9128	107	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 6.

All (107) 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:278:SER:OG	1:C:154:HIS:HD2	1.81	0.64
1:D:94:ARG:HH11	1:D:114:HIS:HD2	1.47	0.62
1:C:94:ARG:HH11	1:C:114:HIS:HD2	1.47	0.62
1:C:15:HIS:HE1	1:C:49:LEU:H	1.47	0.62
1:D:15:HIS:HE1	1:D:49:LEU:H	1.48	0.61
1:B:94:ARG:HH11	1:B:114:HIS:HD2	1.47	0.61
1:B:310:HIS:CD2	1:B:313:ARG:HD2	2.36	0.61
1:B:15:HIS:HE1	1:B:49:LEU:H	1.47	0.60
1:A:94:ARG:HH11	1:A:114:HIS:HD2	1.49	0.60
1:D:310:HIS:CD2	1:D:313:ARG:HD2	2.38	0.58
1:D:310:HIS:HD2	1:D:313:ARG:HH11	1.51	0.58
1:A:15:HIS:HE1	1:A:49:LEU:H	1.50	0.58
1:C:310:HIS:CD2	1:C:313:ARG:HD2	2.38	0.58
1:B:310:HIS:HD2	1:B:313:ARG:HH11	1.51	0.57
1:C:310:HIS:HD2	1:C:313:ARG:HH11	1.51	0.57
1:A:310:HIS:CD2	1:A:313:ARG:HD2	2.39	0.57
1:A:310:HIS:HD2	1:A:313:ARG:HH11	1.51	0.57
1:C:109:THR:HG21	1:C:114:HIS:HE1	1.68	0.57
1:D:109:THR:HG21	1:D:114:HIS:HE1	1.70	0.57
1:B:109:THR:HG21	1:B:114:HIS:HE1	1.70	0.57
1:A:109:THR:HG21	1:A:114:HIS:HE1	1.71	0.56
1:D:192:THR:HG23	1:D:272:LEU:HB3	1.88	0.56
1:C:150:VAL:HG12	1:C:151:SER:H	1.70	0.55
1:A:192:THR:HG23	1:A:272:LEU:HB3	1.88	0.55
1:A:286:LEU:HB3	1:A:291:GLN:HB3	1.89	0.55
1:C:15:HIS:CE1	1:C:49:LEU:H	2.25	0.55
1:C:150:VAL:CG1	1:C:151:SER:H	2.20	0.54
1:C:192:THR:HG23	1:C:272:LEU:HB3	1.89	0.53
1:B:15:HIS:CE1	1:B:49:LEU:H	2.26	0.53
1:B:296:ARG:HB3	1:B:300:ALA:HB3	1.91	0.53
1:B:64:GLN:O	1:B:67:ALA:HB3	2.09	0.53
1:C:286:LEU:HB3	1:C:291:GLN:HB3	1.90	0.53
1:D:15:HIS:CE1	1:D:49:LEU:H	2.27	0.52
1:B:286:LEU:HB3	1:B:291:GLN:HB3	1.90	0.52
1:C:177:ARG:NH1	1:D:301:GLU:OE2	2.42	0.52
1:B:192:THR:HG23	1:B:272:LEU:HB3	1.90	0.52
1:A:296:ARG:HB3	1:A:300:ALA:HB3	1.92	0.51
1:D:286:LEU:HB3	1:D:291:GLN:HB3	1.92	0.50
1:D:310:HIS:O	1:D:313:ARG:HB3	2.12	0.50
1:D:296:ARG:HB3	1:D:300:ALA:HB3	1.91	0.50
1:C:296:ARG:HB3	1:C:300:ALA:HB3	1.92	0.50
1:B:219:THR:HA	1:B:223:VAL:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:113:PRO:HB3	1:D:266:VAL:HG22	1.95	0.49
1:B:310:HIS:O	1:B:313:ARG:HB3	2.12	0.49
1:A:15:HIS:CE1	1:A:49:LEU:H	2.29	0.49
1:C:310:HIS:CD2	1:C:313:ARG:HH11	2.31	0.49
1:B:224:THR:HG21	1:A:191:GLN:HG2	1.95	0.49
1:C:109:THR:HG21	1:C:114:HIS:CE1	2.48	0.48
1:D:310:HIS:CD2	1:D:313:ARG:HH11	2.31	0.48
1:C:113:PRO:HB3	1:C:266:VAL:HG22	1.95	0.48
1:A:310:HIS:O	1:A:313:ARG:HB3	2.14	0.48
1:A:310:HIS:CD2	1:A:313:ARG:HH11	2.30	0.48
1:D:215:GLN:O	1:D:226:ALA:HA	2.14	0.48
1:D:109:THR:HG21	1:D:114:HIS:CE1	2.48	0.47
1:C:310:HIS:O	1:C:313:ARG:HB3	2.14	0.47
1:A:109:THR:HG21	1:A:114:HIS:CE1	2.49	0.47
1:D:155:ASN:O	1:D:157:ASP:N	2.47	0.47
1:A:113:PRO:HB3	1:A:266:VAL:HG22	1.96	0.47
1:B:310:HIS:CD2	1:B:313:ARG:HH11	2.30	0.47
1:B:22:LYS:HD2	1:B:23:PHE:CE2	2.50	0.47
1:C:196:THR:HA	1:C:204:HIS:O	2.15	0.47
1:B:113:PRO:HB3	1:B:266:VAL:HG22	1.96	0.46
1:D:236:VAL:HG12	1:D:240:THR:HG23	1.98	0.46
1:C:121:ASP:O	1:C:124:GLN:HG2	2.16	0.46
1:B:196:THR:HA	1:B:204:HIS:O	2.16	0.46
1:D:196:THR:HA	1:D:204:HIS:O	2.15	0.46
1:D:60:GLY:O	1:D:64:GLN:HG2	2.16	0.45
1:A:196:THR:HA	1:A:204:HIS:O	2.17	0.45
1:D:20:THR:HA	1:D:51:GLY:HA2	1.98	0.45
1:C:60:GLY:O	1:C:64:GLN:HG2	2.15	0.45
1:B:60:GLY:O	1:B:64:GLN:HG2	2.16	0.45
1:B:215:GLN:O	1:B:226:ALA:HA	2.16	0.45
1:A:280:SER:HB2	1:C:154:HIS:CD2	2.51	0.45
1:B:109:THR:HG21	1:B:114:HIS:CE1	2.49	0.44
1:C:26:VAL:HG21	1:C:48:ASN:OD1	2.16	0.44
1:B:123:VAL:O	1:B:127:LEU:HG	2.17	0.44
1:A:114:HIS:CE1	1:A:273:PHE:HB2	2.53	0.44
1:A:121:ASP:O	1:A:124:GLN:HG2	2.17	0.44
1:B:22:LYS:HD2	1:B:23:PHE:CZ	2.52	0.44
1:B:121:ASP:O	1:B:124:GLN:HG2	2.18	0.44
1:A:22:LYS:HD3	1:A:51:GLY:HA3	1.99	0.44
1:D:121:ASP:O	1:D:124:GLN:HG2	2.17	0.44
1:A:215:GLN:O	1:A:226:ALA:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:GLY:O	1:A:64:GLN:HG2	2.17	0.44
1:A:123:VAL:O	1:A:127:LEU:HG	2.17	0.43
1:C:123:VAL:O	1:C:127:LEU:HG	2.19	0.43
1:C:215:GLN:O	1:C:226:ALA:HA	2.19	0.43
1:A:25:ASN:OD1	1:A:27:VAL:HG12	2.19	0.43
1:B:114:HIS:CE1	1:B:273:PHE:HB2	2.54	0.43
1:D:234:LEU:HD23	1:D:234:LEU:HA	1.83	0.42
1:C:114:HIS:CE1	1:C:273:PHE:HB2	2.54	0.42
1:C:150:VAL:CG1	1:C:151:SER:N	2.82	0.42
1:B:310:HIS:HD2	1:B:313:ARG:HD2	1.82	0.42
1:D:24:ALA:O	1:D:26:VAL:HG23	2.20	0.42
1:A:67:ALA:O	1:A:71:GLN:HG3	2.18	0.42
1:D:114:HIS:CE1	1:D:273:PHE:HB2	2.54	0.41
1:C:177:ARG:HH11	1:C:177:ARG:HG2	1.85	0.41
1:D:123:VAL:O	1:D:127:LEU:HG	2.19	0.41
1:C:310:HIS:HD2	1:C:313:ARG:HD2	1.86	0.41
1:D:26:VAL:HG21	1:D:48:ASN:CG	2.41	0.41
1:D:302:ASP:HA	1:D:303:PRO:HD3	1.94	0.41
1:C:45:TYR:HB3	1:C:68:TYR:OH	2.20	0.41
1:B:283:TRP:CH2	1:B:306:VAL:HG21	2.56	0.41
1:C:215:GLN:OE1	1:C:261:GLN:HG2	2.21	0.40
1:D:194:ALA:O	1:D:274:GLY:HA2	2.22	0.40
1:B:45:TYR:HB3	1:B:68:TYR:OH	2.21	0.40
1:D:283:TRP:CH2	1:D:306:VAL:HG21	2.57	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	316/318 (99%)	298 (94%)	14 (4%)	4 (1%)	15 53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	316/318 (99%)	295 (93%)	17 (5%)	4 (1%)	15 53
1	C	316/318 (99%)	298 (94%)	14 (4%)	4 (1%)	15 53
1	D	316/318 (99%)	291 (92%)	17 (5%)	8 (2%)	7 34
All	All	1264/1272 (99%)	1182 (94%)	62 (5%)	20 (2%)	12 48

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	134	LEU
1	B	261	GLN
1	A	261	GLN
1	C	261	GLN
1	D	154	HIS
1	D	156	THR
1	D	218	SER
1	D	261	GLN
1	B	150	VAL
1	B	292	LEU
1	A	134	LEU
1	A	292	LEU
1	C	150	VAL
1	C	151	SER
1	C	292	LEU
1	D	292	LEU
1	A	156	THR
1	D	25	ASN
1	D	152	SER
1	D	151	SER

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	248/248 (100%)	234 (94%)	14 (6%)	26 65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	248/248 (100%)	231 (93%)	17 (7%)	19 56
1	C	248/248 (100%)	230 (93%)	18 (7%)	17 52
1	D	248/248 (100%)	234 (94%)	14 (6%)	26 65
All	All	992/992 (100%)	929 (94%)	63 (6%)	22 59

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	15	HIS
1	B	27	VAL
1	B	66	LEU
1	B	87	SER
1	B	92	THR
1	B	154	HIS
1	B	157	ASP
1	B	158	GLN
1	B	177	ARG
1	B	192	THR
1	B	196	THR
1	B	227	THR
1	B	233	THR
1	B	236	VAL
1	B	239	VAL
1	B	262	ASN
1	B	292	LEU
1	A	15	HIS
1	A	22	LYS
1	A	26	VAL
1	A	66	LEU
1	A	87	SER
1	A	92	THR
1	A	151	SER
1	A	177	ARG
1	A	192	THR
1	A	196	THR
1	A	227	THR
1	A	239	VAL
1	A	262	ASN
1	A	292	LEU
1	C	15	HIS
1	C	27	VAL

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Mol	Chain	Res	Type
1	C	66	LEU
1	C	87	SER
1	C	92	THR
1	C	150	VAL
1	C	152	SER
1	C	153	SER
1	C	177	ARG
1	C	192	THR
1	C	196	THR
1	C	218	SER
1	C	227	THR
1	C	236	VAL
1	C	239	VAL
1	C	262	ASN
1	C	280	SER
1	C	292	LEU
1	D	15	HIS
1	D	20	THR
1	D	27	VAL
1	D	66	LEU
1	D	87	SER
1	D	92	THR
1	D	155	ASN
1	D	177	ARG
1	D	192	THR
1	D	196	THR
1	D	227	THR
1	D	239	VAL
1	D	262	ASN
1	D	292	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	15	HIS
1	B	88	GLN
1	B	114	HIS
1	B	310	HIS
1	A	15	HIS
1	A	88	GLN
1	A	114	HIS
1	A	310	HIS

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Mol	Chain	Res	Type
1	C	15	HIS
1	C	88	GLN
1	C	114	HIS
1	C	154	HIS
1	C	310	HIS
1	D	15	HIS
1	D	88	GLN
1	D	114	HIS
1	D	310	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

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.