



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:04 PM GMT

PDB ID : 4KEA
Title : Crystal structure of D196N mutant of Monoglyceride lipase from Bacillus sp. H257 in space group P212121
Authors : Rengachari, S.; Aschauer, P.; Gruber, K.; Dreveny, I.; Oberer, M.
Deposited on : 2013-04-25
Resolution : 1.70 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:

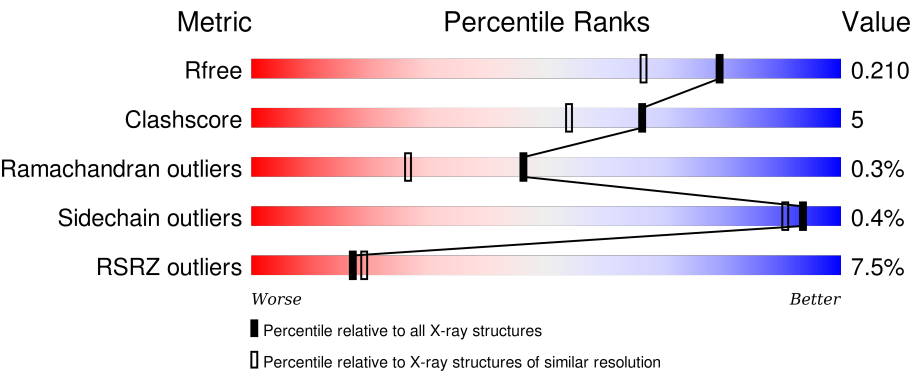
MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : 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 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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 $\leq 5\%$. 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	270	<div><div>4%</div><div>86%6%8%</div></div>
1	B	270	<div><div>4%</div><div>83%7%10%</div></div>
1	C	270	<div><div>7%</div><div>82%9%9%</div></div>
1	D	270	<div><div>7%</div><div>86%10%</div></div>
1	E	270	<div><div>8%</div><div>86%10%</div></div>

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Mol	Chain	Length	Quality of chain
1	F	270	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MPD	A	301	-	-	-	X
2	MPD	A	303	-	-	-	X
2	MPD	A	304	-	-	-	X
2	MPD	A	305	-	-	-	X
2	MPD	A	306	-	-	-	X
2	MPD	B	302	-	-	-	X
2	MPD	B	304	-	-	-	X
2	MPD	C	302	-	-	-	X
2	MPD	C	303	-	-	-	X
2	MPD	C	304	-	-	X	X
2	MPD	E	303	-	-	-	X
2	MPD	E	304	-	-	-	X
2	MPD	F	301	-	-	-	X
2	MPD	F	303	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 12656 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 Thermostable monoacylglycerol lipase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	13	0
			1983	1269	331	371	12			
1	B	243	Total	C	N	O	S	0	7	0
			1921	1232	320	359	10			
1	C	245	Total	C	N	O	S	0	10	0
			1957	1254	328	364	11			
1	D	244	Total	C	N	O	S	0	5	0
			1918	1227	322	359	10			
1	E	242	Total	C	N	O	S	0	3	0
			1896	1211	321	355	9			
1	F	240	Total	C	N	O	S	0	4	0
			1885	1208	316	350	11			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P82597
A	-18	GLY	-	EXPRESSION TAG	UNP P82597
A	-17	SER	-	EXPRESSION TAG	UNP P82597
A	-16	SER	-	EXPRESSION TAG	UNP P82597
A	-15	HIS	-	EXPRESSION TAG	UNP P82597
A	-14	HIS	-	EXPRESSION TAG	UNP P82597
A	-13	HIS	-	EXPRESSION TAG	UNP P82597
A	-12	HIS	-	EXPRESSION TAG	UNP P82597
A	-11	HIS	-	EXPRESSION TAG	UNP P82597
A	-10	HIS	-	EXPRESSION TAG	UNP P82597
A	-9	SER	-	EXPRESSION TAG	UNP P82597
A	-8	SER	-	EXPRESSION TAG	UNP P82597
A	-7	GLY	-	EXPRESSION TAG	UNP P82597
A	-6	LEU	-	EXPRESSION TAG	UNP P82597
A	-5	VAL	-	EXPRESSION TAG	UNP P82597
A	-4	PRO	-	EXPRESSION TAG	UNP P82597
A	-3	ARG	-	EXPRESSION TAG	UNP P82597

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P82597
A	-1	SER	-	EXPRESSION TAG	UNP P82597
A	0	HIS	-	EXPRESSION TAG	UNP P82597
A	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597
B	-19	MET	-	EXPRESSION TAG	UNP P82597
B	-18	GLY	-	EXPRESSION TAG	UNP P82597
B	-17	SER	-	EXPRESSION TAG	UNP P82597
B	-16	SER	-	EXPRESSION TAG	UNP P82597
B	-15	HIS	-	EXPRESSION TAG	UNP P82597
B	-14	HIS	-	EXPRESSION TAG	UNP P82597
B	-13	HIS	-	EXPRESSION TAG	UNP P82597
B	-12	HIS	-	EXPRESSION TAG	UNP P82597
B	-11	HIS	-	EXPRESSION TAG	UNP P82597
B	-10	HIS	-	EXPRESSION TAG	UNP P82597
B	-9	SER	-	EXPRESSION TAG	UNP P82597
B	-8	SER	-	EXPRESSION TAG	UNP P82597
B	-7	GLY	-	EXPRESSION TAG	UNP P82597
B	-6	LEU	-	EXPRESSION TAG	UNP P82597
B	-5	VAL	-	EXPRESSION TAG	UNP P82597
B	-4	PRO	-	EXPRESSION TAG	UNP P82597
B	-3	ARG	-	EXPRESSION TAG	UNP P82597
B	-2	GLY	-	EXPRESSION TAG	UNP P82597
B	-1	SER	-	EXPRESSION TAG	UNP P82597
B	0	HIS	-	EXPRESSION TAG	UNP P82597
B	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597
C	-19	MET	-	EXPRESSION TAG	UNP P82597
C	-18	GLY	-	EXPRESSION TAG	UNP P82597
C	-17	SER	-	EXPRESSION TAG	UNP P82597
C	-16	SER	-	EXPRESSION TAG	UNP P82597
C	-15	HIS	-	EXPRESSION TAG	UNP P82597
C	-14	HIS	-	EXPRESSION TAG	UNP P82597
C	-13	HIS	-	EXPRESSION TAG	UNP P82597
C	-12	HIS	-	EXPRESSION TAG	UNP P82597
C	-11	HIS	-	EXPRESSION TAG	UNP P82597
C	-10	HIS	-	EXPRESSION TAG	UNP P82597
C	-9	SER	-	EXPRESSION TAG	UNP P82597
C	-8	SER	-	EXPRESSION TAG	UNP P82597
C	-7	GLY	-	EXPRESSION TAG	UNP P82597
C	-6	LEU	-	EXPRESSION TAG	UNP P82597
C	-5	VAL	-	EXPRESSION TAG	UNP P82597
C	-4	PRO	-	EXPRESSION TAG	UNP P82597
C	-3	ARG	-	EXPRESSION TAG	UNP P82597

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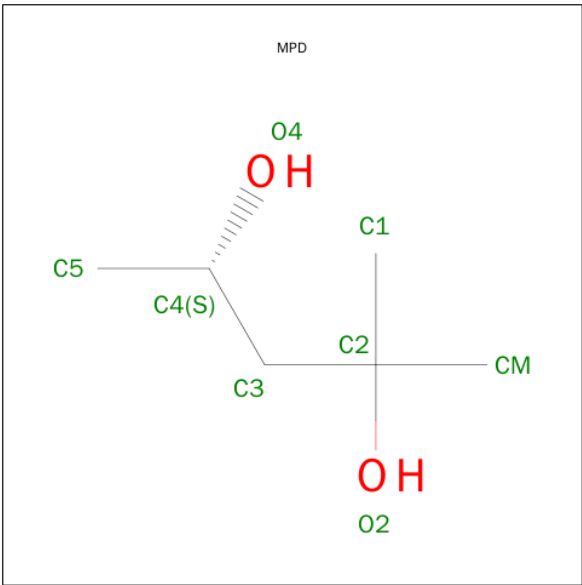
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	EXPRESSION TAG	UNP P82597
C	-1	SER	-	EXPRESSION TAG	UNP P82597
C	0	HIS	-	EXPRESSION TAG	UNP P82597
C	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597
D	-19	MET	-	EXPRESSION TAG	UNP P82597
D	-18	GLY	-	EXPRESSION TAG	UNP P82597
D	-17	SER	-	EXPRESSION TAG	UNP P82597
D	-16	SER	-	EXPRESSION TAG	UNP P82597
D	-15	HIS	-	EXPRESSION TAG	UNP P82597
D	-14	HIS	-	EXPRESSION TAG	UNP P82597
D	-13	HIS	-	EXPRESSION TAG	UNP P82597
D	-12	HIS	-	EXPRESSION TAG	UNP P82597
D	-11	HIS	-	EXPRESSION TAG	UNP P82597
D	-10	HIS	-	EXPRESSION TAG	UNP P82597
D	-9	SER	-	EXPRESSION TAG	UNP P82597
D	-8	SER	-	EXPRESSION TAG	UNP P82597
D	-7	GLY	-	EXPRESSION TAG	UNP P82597
D	-6	LEU	-	EXPRESSION TAG	UNP P82597
D	-5	VAL	-	EXPRESSION TAG	UNP P82597
D	-4	PRO	-	EXPRESSION TAG	UNP P82597
D	-3	ARG	-	EXPRESSION TAG	UNP P82597
D	-2	GLY	-	EXPRESSION TAG	UNP P82597
D	-1	SER	-	EXPRESSION TAG	UNP P82597
D	0	HIS	-	EXPRESSION TAG	UNP P82597
D	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597
E	-19	MET	-	EXPRESSION TAG	UNP P82597
E	-18	GLY	-	EXPRESSION TAG	UNP P82597
E	-17	SER	-	EXPRESSION TAG	UNP P82597
E	-16	SER	-	EXPRESSION TAG	UNP P82597
E	-15	HIS	-	EXPRESSION TAG	UNP P82597
E	-14	HIS	-	EXPRESSION TAG	UNP P82597
E	-13	HIS	-	EXPRESSION TAG	UNP P82597
E	-12	HIS	-	EXPRESSION TAG	UNP P82597
E	-11	HIS	-	EXPRESSION TAG	UNP P82597
E	-10	HIS	-	EXPRESSION TAG	UNP P82597
E	-9	SER	-	EXPRESSION TAG	UNP P82597
E	-8	SER	-	EXPRESSION TAG	UNP P82597
E	-7	GLY	-	EXPRESSION TAG	UNP P82597
E	-6	LEU	-	EXPRESSION TAG	UNP P82597
E	-5	VAL	-	EXPRESSION TAG	UNP P82597
E	-4	PRO	-	EXPRESSION TAG	UNP P82597
E	-3	ARG	-	EXPRESSION TAG	UNP P82597

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP P82597
E	-1	SER	-	EXPRESSION TAG	UNP P82597
E	0	HIS	-	EXPRESSION TAG	UNP P82597
E	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597
F	-19	MET	-	EXPRESSION TAG	UNP P82597
F	-18	GLY	-	EXPRESSION TAG	UNP P82597
F	-17	SER	-	EXPRESSION TAG	UNP P82597
F	-16	SER	-	EXPRESSION TAG	UNP P82597
F	-15	HIS	-	EXPRESSION TAG	UNP P82597
F	-14	HIS	-	EXPRESSION TAG	UNP P82597
F	-13	HIS	-	EXPRESSION TAG	UNP P82597
F	-12	HIS	-	EXPRESSION TAG	UNP P82597
F	-11	HIS	-	EXPRESSION TAG	UNP P82597
F	-10	HIS	-	EXPRESSION TAG	UNP P82597
F	-9	SER	-	EXPRESSION TAG	UNP P82597
F	-8	SER	-	EXPRESSION TAG	UNP P82597
F	-7	GLY	-	EXPRESSION TAG	UNP P82597
F	-6	LEU	-	EXPRESSION TAG	UNP P82597
F	-5	VAL	-	EXPRESSION TAG	UNP P82597
F	-4	PRO	-	EXPRESSION TAG	UNP P82597
F	-3	ARG	-	EXPRESSION TAG	UNP P82597
F	-2	GLY	-	EXPRESSION TAG	UNP P82597
F	-1	SER	-	EXPRESSION TAG	UNP P82597
F	0	HIS	-	EXPRESSION TAG	UNP P82597
F	196	ASN	ASP	ENGINEERED MUTATION	UNP P82597

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			8	6	2		
2	D	1	Total	C	O	0	0
			8	6	2		
2	D	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	F	1	Total	C	O	0	0
			8	6	2		
2	F	1	Total	C	O	0	0
			8	6	2		
2	F	1	Total	C	O	0	0
			8	6	2		

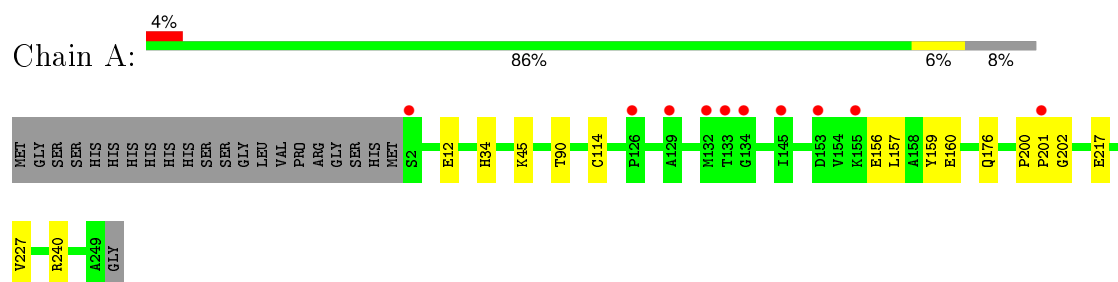
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	207	Total	O	0	0
			207	207		
3	B	188	Total	O	0	0
			188	188		
3	C	123	Total	O	0	0
			123	123		
3	D	125	Total	O	0	0
			125	125		
3	E	153	Total	O	0	0
			153	153		
3	F	108	Total	O	0	0
			108	108		

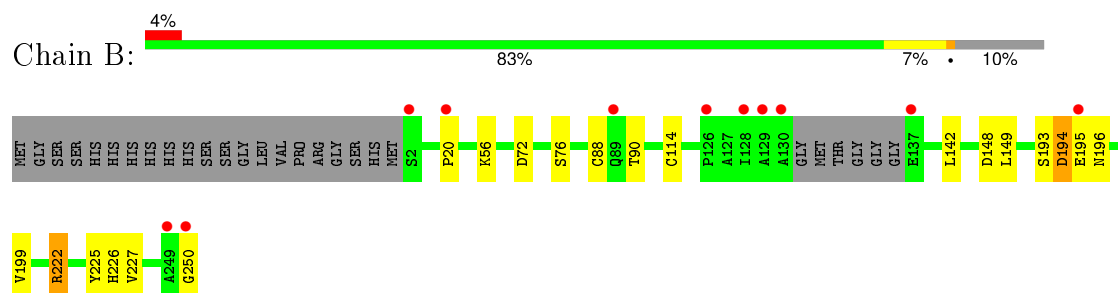
3 Residue-property plots [i](#)

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.

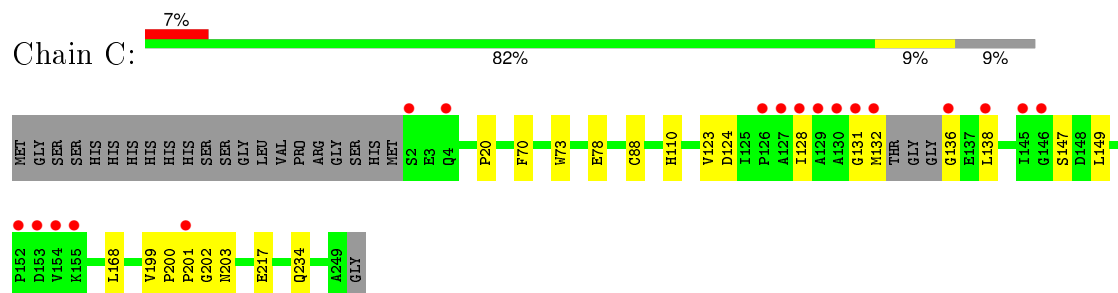
- Molecule 1: Thermostable monoacylglycerol lipase



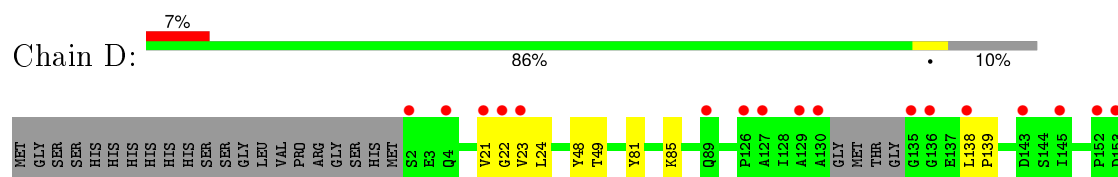
- Molecule 1: Thermostable monoacylglycerol lipase

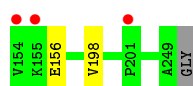


- Molecule 1: Thermostable monoacylglycerol lipase

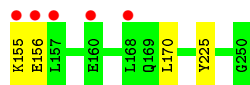
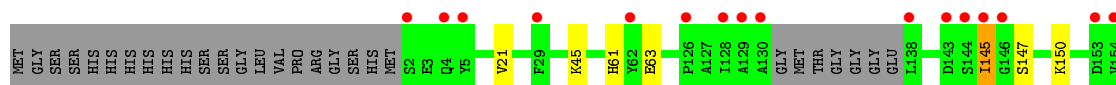
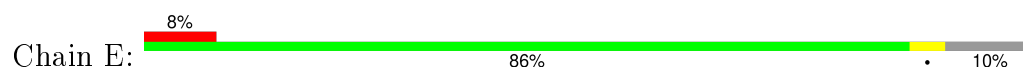


- Molecule 1: Thermostable monoacylglycerol lipase

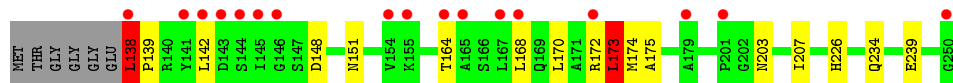
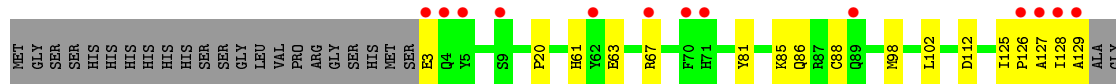
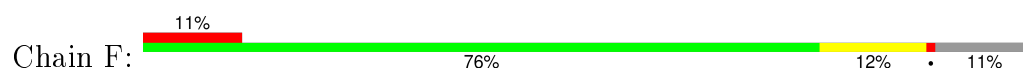




- Molecule 1: Thermostable monoacylglycerol lipase



- Molecule 1: Thermostable monoacylglycerol lipase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.16Å 183.13Å 244.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.88 – 1.70 19.88 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.88-1.70) 98.9 (19.88-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.04 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.191 , 0.211 0.191 , 0.210	Depositor DCC
R_{free} test set	9723 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 195083 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12656	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.44	2/2057 (0.1%)	0.47	0/2799
1	B	0.30	0/1986	0.46	0/2701
1	C	0.37	0/2021	0.45	0/2750
1	D	0.29	0/1977	0.43	0/2688
1	E	0.31	0/1949	0.45	0/2651
1	F	0.35	0/1941	0.49	1/2640 (0.0%)
All	All	0.35	2/11931 (0.0%)	0.46	1/16229 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	159	TYR	CG-CD1	-5.64	1.31	1.39
1	A	12	GLU	CD-OE1	-5.29	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	138	LEU	CA-CB-CG	5.10	127.04	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1983	0	1978	15	0
1	B	1921	0	1916	19	0
1	C	1957	0	1947	22	0
1	D	1918	0	1902	8	0
1	E	1896	0	1874	10	0
1	F	1885	0	1872	33	0
2	A	48	0	83	6	0
2	B	32	0	56	4	0
2	C	32	0	55	7	0
2	D	24	0	42	2	0
2	E	32	0	56	5	0
2	F	24	0	42	3	0
3	A	207	0	0	4	0
3	B	188	0	0	0	0
3	C	123	0	0	3	0
3	D	125	0	0	0	0
3	E	153	0	0	0	0
3	F	108	0	0	4	0
All	All	12656	0	11823	116	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 5.

All (116) 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)
2:C:304:MPD:C5	2:C:304:MPD:H12	1.62	1.25
2:C:304:MPD:H52	2:C:304:MPD:C1	1.69	1.14
1:B:222:ARG:HH11	1:B:222:ARG:HG2	1.27	0.99
2:C:304:MPD:C5	2:C:304:MPD:C1	2.30	0.97
1:C:234:GLN:HB3	2:C:303:MPD:H51	1.57	0.85
1:F:170:LEU:HD13	1:F:170:LEU:C	1.98	0.84
1:E:156:GLU:OE2	2:E:302:MPD:O4	1.96	0.83
2:C:304:MPD:H52	2:C:304:MPD:H12	0.82	0.79
1:F:172:ARG:O	1:F:173:LEU:C	2.25	0.73
1:C:124:ASP:H	1:C:203[B]:ASN:HD21	1.33	0.73
1:B:149[B]:LEU:HD23	1:B:227:VAL:HB	1.70	0.73
1:A:34:HIS:HD2	1:A:156[B]:GLU:OE2	1.71	0.73
1:C:123:VAL:H	1:C:203[B]:ASN:ND2	1.86	0.73
1:F:81:TYR:CE2	1:F:85:LYS:HD2	2.24	0.72
1:C:123:VAL:H	1:C:203[B]:ASN:HD22	1.35	0.72
1:A:217:GLU:HB3	1:C:217:GLU:HB3	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:98:MET:SD	1:F:170:LEU:HD23	2.33	0.68
1:C:200[B]:PRO:HB3	1:C:201[B]:PRO:HD2	1.77	0.66
1:B:222:ARG:NH1	1:B:222:ARG:HG2	2.07	0.66
1:F:170:LEU:HD11	1:F:174:MET:SD	2.38	0.64
2:C:304:MPD:HM1	3:C:519:HOH:O	1.98	0.64
1:A:45[B]:LYS:CD	3:A:605:HOH:O	2.45	0.64
1:F:148:ASP:HB3	1:F:226:HIS:HB3	1.80	0.64
1:D:22:GLY:HA2	1:D:49:THR:O	1.98	0.63
2:B:304:MPD:O4	2:B:304:MPD:O2	2.11	0.63
1:B:250:GLY:OXT	1:E:45:LYS:NZ	2.31	0.62
1:B:193:SER:O	1:B:195:GLU:N	2.33	0.62
1:A:45[B]:LYS:HD3	3:A:605:HOH:O	2.00	0.62
1:F:3:GLU:OE2	1:F:67:ARG:NH2	2.34	0.60
1:F:170:LEU:CD1	1:F:170:LEU:C	2.70	0.60
1:E:21:VAL:HG21	2:E:303:MPD:H12	1.83	0.59
1:F:85:LYS:NZ	1:F:112:ASP:OD2	2.33	0.59
1:F:172:ARG:O	1:F:175:ALA:N	2.35	0.58
2:C:304:MPD:H53	2:C:304:MPD:C1	2.31	0.57
1:B:195:GLU:O	1:B:225:TYR:HA	2.04	0.57
1:D:21:VAL:O	1:D:48:TYR:HD1	1.88	0.56
2:B:303:MPD:HM1	2:B:304:MPD:H51	1.89	0.55
1:E:170:LEU:HD21	2:E:301:MPD:H11	1.89	0.55
1:F:170:LEU:HD13	1:F:170:LEU:O	2.06	0.54
1:F:139:PRO:O	1:F:164:THR:HG21	2.07	0.54
2:A:304:MPD:O2	2:A:304:MPD:H52	2.07	0.54
1:A:90:THR:HG23	1:A:114[B]:CYS:SG	2.48	0.54
1:B:193:SER:C	1:B:195:GLU:H	2.12	0.53
1:C:78:GLU:OE2	1:C:110:HIS:NE2	2.25	0.53
1:E:147:SER:OG	1:E:156:GLU:OE1	2.18	0.53
1:C:136:GLY:N	3:C:484:HOH:O	2.41	0.53
1:B:193:SER:C	1:B:195:GLU:N	2.62	0.53
1:C:124:ASP:H	1:C:203[B]:ASN:ND2	2.04	0.52
1:B:56[B]:LYS:HD2	1:B:72:ASP:HB3	1.92	0.52
2:F:303:MPD:O2	2:F:303:MPD:O4	2.16	0.52
1:E:156:GLU:CD	2:E:302:MPD:HO4	2.04	0.51
1:B:90:THR:HG23	1:B:114[B]:CYS:SG	2.51	0.51
1:F:20:PRO:O	1:F:88:CYS:HB3	2.11	0.51
1:F:168:LEU:O	1:F:172:ARG:HG3	2.10	0.50
1:E:150:LYS:HE3	1:E:225:TYR:CE1	2.45	0.50
1:B:20:PRO:O	1:B:88:CYS:HB3	2.11	0.50
1:D:198:VAL:HG13	2:D:302:MPD:H11	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ARG:HH12	2:A:303:MPD:H12	1.77	0.50
1:C:147:SER:HB3	1:C:149:LEU:HD12	1.95	0.49
1:F:3:GLU:CD	3:F:508:HOH:O	2.51	0.49
1:F:234:GLN:HB3	2:F:303:MPD:HM2	1.95	0.49
1:F:3:GLU:CG	3:F:508:HOH:O	2.60	0.49
1:F:172:ARG:O	1:F:174:MET:N	2.46	0.48
1:B:56[A]:LYS:HG3	1:B:76:SER:OG	2.12	0.48
1:F:170:LEU:CD1	1:F:174:MET:SD	3.02	0.47
1:F:203:ASN:O	1:F:207[A]:ILE:HG12	2.14	0.47
1:F:142:LEU:CD2	1:F:164:THR:HG22	2.44	0.47
1:B:56[A]:LYS:HD2	1:B:72:ASP:HB3	1.96	0.47
1:A:176:GLN:NE2	3:A:517:HOH:O	2.33	0.46
1:F:138:LEU:HB2	1:F:139:PRO:HD2	1.98	0.46
2:E:302:MPD:O2	2:E:302:MPD:O4	2.20	0.46
1:F:142:LEU:HD22	1:F:164:THR:HG22	1.98	0.46
1:B:222:ARG:CG	1:B:222:ARG:HH11	2.06	0.46
1:D:81:TYR:CZ	1:D:85:LYS:HD2	2.51	0.46
1:F:61:HIS:CE1	1:F:63:GLU:HG3	2.50	0.46
1:A:157:LEU:HG	2:A:306:MPD:H4	1.98	0.45
1:F:151:ASN:HD22	2:F:303:MPD:C5	2.29	0.45
2:A:306:MPD:O2	2:A:306:MPD:O4	2.26	0.45
1:E:61:HIS:CE1	1:E:63:GLU:HG3	2.52	0.45
1:A:45[B]:LYS:NZ	3:A:605:HOH:O	2.49	0.45
1:A:217:GLU:CB	1:C:217:GLU:HB3	2.44	0.45
1:C:132:MET:HE1	1:C:168[B]:LEU:HB2	1.98	0.44
1:C:20:PRO:O	1:C:88:CYS:HB3	2.17	0.44
1:C:128:ILE:O	1:C:132:MET:HG3	2.18	0.44
1:E:145:ILE:HG13	1:E:145:ILE:H	1.60	0.44
1:F:148:ASP:HB3	1:F:226:HIS:CB	2.46	0.44
1:C:200[B]:PRO:CB	1:C:201[B]:PRO:HD2	2.30	0.43
1:C:132:MET:HE1	1:C:168[A]:LEU:HB2	2.00	0.43
1:A:227:VAL:HG22	2:A:301:MPD:H53	2.00	0.43
1:A:200[B]:PRO:CB	1:A:201[B]:PRO:HD2	2.47	0.43
1:B:142:LEU:HB3	2:B:304:MPD:H53	2.01	0.43
1:B:196:ASN:HB3	1:B:199:VAL:O	2.18	0.43
1:D:22:GLY:CA	1:D:49:THR:O	2.66	0.43
1:C:201[B]:PRO:HA	1:C:202[B]:GLY:HA2	1.83	0.43
1:A:201[B]:PRO:HA	1:A:202[B]:GLY:HA2	1.59	0.43
1:F:102:LEU:HD21	1:F:173:LEU:HD21	2.01	0.42
1:E:155:LYS:HD2	1:E:155:LYS:HA	1.81	0.42
1:B:222:ARG:CG	1:B:222:ARG:NH1	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:148:ASP:HB3	1:B:226:HIS:HB3	2.02	0.41
1:C:70:PHE:HA	1:C:73:TRP:CE3	2.55	0.41
1:F:86:GLN:NE2	3:F:498:HOH:O	2.46	0.41
1:A:160:GLU:HG3	2:A:306:MPD:C1	2.51	0.41
1:D:138:LEU:HA	1:D:139:PRO:HD3	1.93	0.41
1:D:156:GLU:OE2	2:D:301:MPD:O4	2.28	0.41
1:C:201[B]:PRO:O	1:C:201[B]:PRO:HG2	2.20	0.41
1:F:128:ILE:HG13	1:F:129:ALA:N	2.35	0.41
1:F:172:ARG:C	1:F:174:MET:N	2.71	0.41
1:C:199[B]:VAL:HA	1:C:200[B]:PRO:HD3	1.94	0.41
1:D:23:VAL:HG12	1:D:24:LEU:N	2.35	0.41
1:F:239:GLU:OE2	3:F:505:HOH:O	2.22	0.41
1:C:217:GLU:OE1	3:C:463:HOH:O	2.22	0.41
1:B:142:LEU:HD13	2:B:304:MPD:H32	2.04	0.40
1:C:138:LEU:HD23	1:C:138:LEU:HA	1.93	0.40
1:F:125:ILE:HA	1:F:126:PRO:HD3	1.80	0.40
1:F:173:LEU:HD12	1:F:173:LEU:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	259/270 (96%)	250 (96%)	9 (4%)	0	100	100
1	B	246/270 (91%)	238 (97%)	7 (3%)	1 (0%)	39	20
1	C	251/270 (93%)	238 (95%)	12 (5%)	1 (0%)	39	20
1	D	245/270 (91%)	237 (97%)	8 (3%)	0	100	100
1	E	241/270 (89%)	232 (96%)	9 (4%)	0	100	100
1	F	240/270 (89%)	228 (95%)	10 (4%)	2 (1%)	24	7
All	All	1482/1620 (92%)	1423 (96%)	55 (4%)	4 (0%)	46	26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	194	ASP
1	C	131	GLY
1	F	127	ALA
1	F	173	LEU

5.3.2 Protein sidechains ⓘ

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	216/222 (97%)	216 (100%)	0	100	100
1	B	209/222 (94%)	207 (99%)	2 (1%)	82	72
1	C	212/222 (96%)	212 (100%)	0	100	100
1	D	207/222 (93%)	207 (100%)	0	100	100
1	E	204/222 (92%)	203 (100%)	1 (0%)	92	88
1	F	204/222 (92%)	202 (99%)	2 (1%)	82	72
All	All	1252/1332 (94%)	1247 (100%)	5 (0%)	93	90

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	194	ASP
1	B	222	ARG
1	E	145	ILE
1	F	138	LEU
1	F	173	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	176	GLN

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 ⓘ

24 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MPD	A	301	-	6,7,7	0.29	0	7,10,10	0.51	0
2	MPD	A	302	-	6,7,7	0.25	0	7,10,10	0.39	0
2	MPD	A	303	-	6,7,7	0.28	0	7,10,10	0.27	0
2	MPD	A	304	-	6,7,7	1.45	1 (16%)	7,10,10	0.62	0
2	MPD	A	305	-	6,7,7	0.23	0	7,10,10	0.41	0
2	MPD	A	306	-	6,7,7	0.26	0	7,10,10	0.21	0
2	MPD	B	301	-	6,7,7	0.26	0	7,10,10	0.36	0
2	MPD	B	302	-	6,7,7	0.26	0	7,10,10	0.31	0
2	MPD	B	303	-	6,7,7	0.20	0	7,10,10	0.43	0
2	MPD	B	304	-	6,7,7	0.22	0	7,10,10	0.28	0
2	MPD	C	301	-	6,7,7	0.24	0	7,10,10	0.43	0
2	MPD	C	302	-	6,7,7	1.00	0	7,10,10	0.96	0
2	MPD	C	303	-	6,7,7	0.26	0	7,10,10	0.51	0
2	MPD	C	304	-	6,7,7	0.64	0	7,10,10	0.79	0
2	MPD	D	301	-	6,7,7	0.32	0	7,10,10	0.36	0
2	MPD	D	302	-	6,7,7	0.26	0	7,10,10	0.33	0
2	MPD	D	303	-	6,7,7	0.25	0	7,10,10	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MPD	E	301	-	6,7,7	0.27	0	7,10,10	0.52	0
2	MPD	E	302	-	6,7,7	0.27	0	7,10,10	0.39	0
2	MPD	E	303	-	6,7,7	0.26	0	7,10,10	0.35	0
2	MPD	E	304	-	6,7,7	0.25	0	7,10,10	0.36	0
2	MPD	F	301	-	6,7,7	0.24	0	7,10,10	0.42	0
2	MPD	F	302	-	6,7,7	0.31	0	7,10,10	0.56	0
2	MPD	F	303	-	6,7,7	0.23	0	7,10,10	0.28	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	A	301	-	-	0/5/5/5	0/0/0/0
2	MPD	A	302	-	-	0/5/5/5	0/0/0/0
2	MPD	A	303	-	-	0/5/5/5	0/0/0/0
2	MPD	A	304	-	-	0/5/5/5	0/0/0/0
2	MPD	A	305	-	-	0/5/5/5	0/0/0/0
2	MPD	A	306	-	-	0/5/5/5	0/0/0/0
2	MPD	B	301	-	-	0/5/5/5	0/0/0/0
2	MPD	B	302	-	-	0/5/5/5	0/0/0/0
2	MPD	B	303	-	-	0/5/5/5	0/0/0/0
2	MPD	B	304	-	-	0/5/5/5	0/0/0/0
2	MPD	C	301	-	-	0/5/5/5	0/0/0/0
2	MPD	C	302	-	-	0/5/5/5	0/0/0/0
2	MPD	C	303	-	-	0/5/5/5	0/0/0/0
2	MPD	C	304	-	-	0/5/5/5	0/0/0/0
2	MPD	D	301	-	-	0/5/5/5	0/0/0/0
2	MPD	D	302	-	-	0/5/5/5	0/0/0/0
2	MPD	D	303	-	-	0/5/5/5	0/0/0/0
2	MPD	E	301	-	-	0/5/5/5	0/0/0/0
2	MPD	E	302	-	-	0/5/5/5	0/0/0/0
2	MPD	E	303	-	-	0/5/5/5	0/0/0/0
2	MPD	E	304	-	-	0/5/5/5	0/0/0/0
2	MPD	F	301	-	-	0/5/5/5	0/0/0/0
2	MPD	F	302	-	-	0/5/5/5	0/0/0/0
2	MPD	F	303	-	-	0/5/5/5	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	304	MPD	O2-C2	-3.09	1.36	1.44

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

14 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	MPD	1	0
2	A	303	MPD	1	0
2	A	304	MPD	1	0
2	A	306	MPD	3	0
2	B	303	MPD	1	0
2	B	304	MPD	4	0
2	C	303	MPD	1	0
2	C	304	MPD	6	0
2	D	301	MPD	1	0
2	D	302	MPD	1	0
2	E	301	MPD	1	0
2	E	302	MPD	3	0
2	E	303	MPD	1	0
2	F	303	MPD	3	0

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	248/270 (91%)	-0.01	10 (4%)	42	46	10, 17, 35, 51	0
1	B	243/270 (90%)	0.12	11 (4%)	37	41	11, 19, 41, 72	0
1	C	245/270 (90%)	0.38	18 (7%)	18	20	12, 25, 51, 81	0
1	D	244/270 (90%)	0.34	20 (8%)	14	16	13, 25, 51, 66	0
1	E	242/270 (89%)	0.40	21 (8%)	13	14	15, 25, 59, 85	0
1	F	240/270 (88%)	0.85	30 (12%)	5	6	17, 33, 61, 83	0
All	All	1462/1620 (90%)	0.34	110 (7%)	17	19	10, 24, 53, 85	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	250	GLY	11.5
1	E	129	ALA	9.4
1	E	138	LEU	9.1
1	F	145	ILE	8.3
1	B	129	ALA	8.1
1	C	131	GLY	6.7
1	B	130	ALA	6.5
1	D	145	ILE	6.4
1	F	138	LEU	6.4
1	F	129	ALA	5.8
1	D	129	ALA	5.7
1	C	130	ALA	5.5
1	B	126	PRO	5.5
1	D	136	GLY	5.4
1	E	145	ILE	5.2
1	E	155	LYS	5.1
1	E	130	ALA	5.0
1	F	146	GLY	4.9
1	C	145	ILE	4.8

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Mol	Chain	Res	Type	RSRZ
1	F	128	ILE	4.8
1	D	23	VAL	4.5
1	F	127	ALA	4.3
1	A	145	ILE	4.3
1	D	138	LEU	4.1
1	D	130	ALA	4.1
1	C	138	LEU	4.1
1	D	2	SER	4.1
1	F	126	PRO	4.0
1	D	135	GLY	4.0
1	A	133	THR	3.9
1	F	141	TYR	3.8
1	E	2	SER	3.8
1	D	155	LYS	3.8
1	C	2	SER	3.8
1	F	155	LYS	3.7
1	E	154	VAL	3.7
1	C	132	MET	3.7
1	E	156	GLU	3.6
1	A	134	GLY	3.6
1	C	126	PRO	3.5
1	E	157	LEU	3.5
1	D	4	GLN	3.4
1	A	129	ALA	3.4
1	A	155	LYS	3.4
1	E	4	GLN	3.4
1	C	155	LYS	3.4
1	F	62	TYR	3.3
1	E	146	GLY	3.3
1	C	154	VAL	3.3
1	D	201	PRO	3.3
1	E	126	PRO	3.1
1	F	5	TYR	3.1
1	C	136	GLY	3.0
1	F	172	ARG	3.0
1	F	168	LEU	3.0
1	C	146	GLY	3.0
1	F	201	PRO	2.9
1	B	89	GLN	2.9
1	F	4	GLN	2.8
1	D	143	ASP	2.8
1	F	71	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	179	ALA	2.7
1	D	154	VAL	2.7
1	B	2	SER	2.7
1	F	144	SER	2.7
1	C	4	GLN	2.7
1	F	142	LEU	2.7
1	A	2	SER	2.7
1	E	128	ILE	2.7
1	C	153[A]	ASP	2.6
1	E	5	TYR	2.6
1	D	153	ASP	2.6
1	E	143	ASP	2.6
1	F	143	ASP	2.6
1	A	153	ASP	2.6
1	C	129	ALA	2.5
1	D	21	VAL	2.5
1	F	3	GLU	2.5
1	B	249	ALA	2.5
1	E	62	TYR	2.4
1	F	9	SER	2.4
1	D	127	ALA	2.4
1	B	128	ILE	2.4
1	A	201[A]	PRO	2.4
1	F	165	ALA	2.3
1	A	126	PRO	2.3
1	B	195	GLU	2.3
1	D	22	GLY	2.3
1	F	89	GLN	2.3
1	A	132	MET	2.3
1	E	153	ASP	2.2
1	F	70	PHE	2.2
1	D	152	PRO	2.2
1	E	168	LEU	2.2
1	C	128	ILE	2.2
1	C	152	PRO	2.1
1	E	29	PHE	2.1
1	B	20	PRO	2.1
1	F	154	VAL	2.1
1	E	144	SER	2.1
1	F	250	GLY	2.1
1	D	89	GLN	2.1
1	F	164	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	127	ALA	2.1
1	F	67	ARG	2.1
1	F	167	LEU	2.1
1	E	160	GLU	2.0
1	B	137	GLU	2.0
1	C	201[A]	PRO	2.0
1	D	126	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 [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MPD	C	303	8/8	0.75	0.25	11.12	44,56,73,74	0
2	MPD	F	303	8/8	0.63	0.43	10.08	28,54,63,65	0
2	MPD	A	306	8/8	0.78	0.27	7.24	39,45,49,49	0
2	MPD	C	304	8/8	0.69	0.29	7.15	55,64,68,71	0
2	MPD	A	305	8/8	0.70	0.30	7.00	35,46,52,55	0
2	MPD	A	304	8/8	0.85	0.19	6.32	32,43,45,47	0
2	MPD	A	303	8/8	0.52	0.28	5.96	39,47,59,59	0
2	MPD	E	304	8/8	0.69	0.22	4.25	43,49,52,52	0
2	MPD	E	303	8/8	0.86	0.18	3.28	19,24,33,35	0
2	MPD	F	301	8/8	0.84	0.20	3.12	47,53,63,70	0
2	MPD	A	301	8/8	0.88	0.17	3.07	31,37,41,42	0
2	MPD	B	304	8/8	0.69	0.20	2.93	33,43,49,49	0
2	MPD	B	302	8/8	0.86	0.16	2.12	19,33,38,42	0
2	MPD	C	302	8/8	0.88	0.17	2.11	39,43,47,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MPD	A	302	8/8	0.91	0.13	1.83	29,33,37,42	0
2	MPD	D	303	8/8	0.80	0.15	1.68	38,44,51,55	0
2	MPD	E	301	8/8	0.88	0.21	1.18	44,52,55,55	0
2	MPD	B	301	8/8	0.91	0.11	0.84	30,36,39,40	0
2	MPD	D	301	8/8	0.84	0.18	0.80	35,47,48,54	0
2	MPD	B	303	8/8	0.93	0.13	0.59	41,44,45,47	0
2	MPD	E	302	8/8	0.85	0.16	0.27	45,54,60,65	0
2	MPD	C	301	8/8	0.90	0.13	0.15	39,43,50,54	0
2	MPD	D	302	8/8	0.88	0.14	0.10	37,47,51,59	0
2	MPD	F	302	8/8	0.92	0.14	-0.19	48,53,56,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.