



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:25 AM UTC

PDB ID : 5LK6 / pdb_00005lk6
Title : Crystal structure of a lipase carboxylesterase from *Sulfolobus islandicus*
Authors : Schwarz-Linnet, T.; Teilum, K.; Olsen, J.G.
Deposited on : 2016-07-21
Resolution : 2.60 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

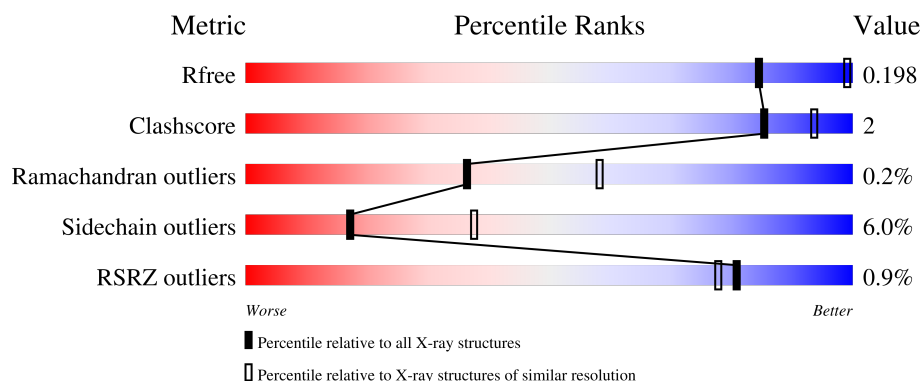
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 of 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	314	% 86% 9% ..
1	B	314	% 85% 8% ..
1	C	314	% 87% 6% ...
1	D	314	85% 9% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9846 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 Alpha/beta hydrolase fold-3 domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2342	1514	391	432	5			
1	B	302	Total	C	N	O	S	0	0	0
			2342	1514	391	432	5			
1	C	302	Total	C	N	O	S	0	0	0
			2342	1514	391	432	5			
1	D	302	Total	C	N	O	S	0	0	0
			2342	1514	391	432	5			

There are 36 discrepancies between the modelled and reference sequences:

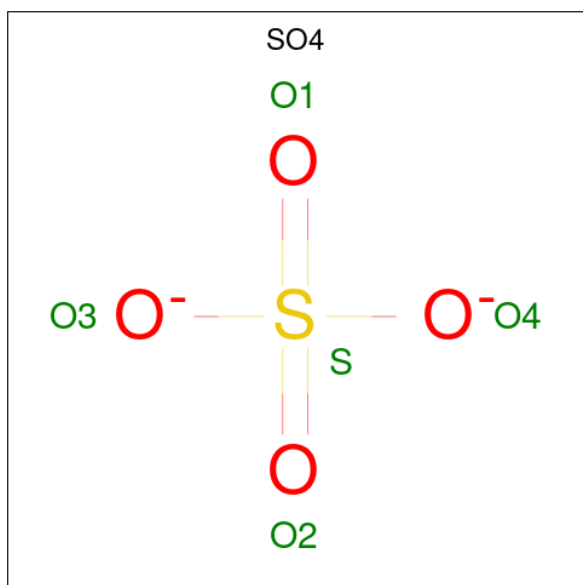
Chain	Residue	Modelled	Actual	Comment	Reference
A	306	ALA	-	expression tag	UNP F0NDQ1
A	307	ALA	-	expression tag	UNP F0NDQ1
A	308	ALA	-	expression tag	UNP F0NDQ1
A	309	HIS	-	expression tag	UNP F0NDQ1
A	310	HIS	-	expression tag	UNP F0NDQ1
A	311	HIS	-	expression tag	UNP F0NDQ1
A	312	HIS	-	expression tag	UNP F0NDQ1
A	313	HIS	-	expression tag	UNP F0NDQ1
A	314	HIS	-	expression tag	UNP F0NDQ1
B	306	ALA	-	expression tag	UNP F0NDQ1
B	307	ALA	-	expression tag	UNP F0NDQ1
B	308	ALA	-	expression tag	UNP F0NDQ1
B	309	HIS	-	expression tag	UNP F0NDQ1
B	310	HIS	-	expression tag	UNP F0NDQ1
B	311	HIS	-	expression tag	UNP F0NDQ1
B	312	HIS	-	expression tag	UNP F0NDQ1
B	313	HIS	-	expression tag	UNP F0NDQ1
B	314	HIS	-	expression tag	UNP F0NDQ1
C	306	ALA	-	expression tag	UNP F0NDQ1
C	307	ALA	-	expression tag	UNP F0NDQ1
C	308	ALA	-	expression tag	UNP F0NDQ1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	309	HIS	-	expression tag	UNP F0NDQ1
C	310	HIS	-	expression tag	UNP F0NDQ1
C	311	HIS	-	expression tag	UNP F0NDQ1
C	312	HIS	-	expression tag	UNP F0NDQ1
C	313	HIS	-	expression tag	UNP F0NDQ1
C	314	HIS	-	expression tag	UNP F0NDQ1
D	306	ALA	-	expression tag	UNP F0NDQ1
D	307	ALA	-	expression tag	UNP F0NDQ1
D	308	ALA	-	expression tag	UNP F0NDQ1
D	309	HIS	-	expression tag	UNP F0NDQ1
D	310	HIS	-	expression tag	UNP F0NDQ1
D	311	HIS	-	expression tag	UNP F0NDQ1
D	312	HIS	-	expression tag	UNP F0NDQ1
D	313	HIS	-	expression tag	UNP F0NDQ1
D	314	HIS	-	expression tag	UNP F0NDQ1

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

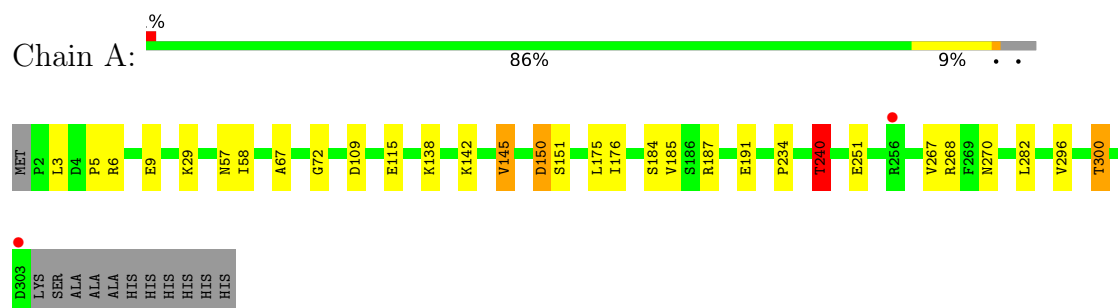
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	104	Total	O	0	0
			104	104		
3	B	118	Total	O	0	0
			118	118		
3	C	109	Total	O	0	0
			109	109		
3	D	107	Total	O	0	0
			107	107		

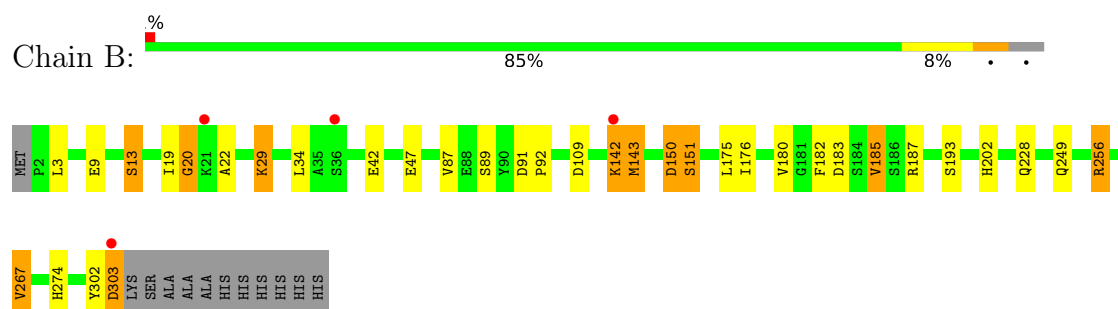
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

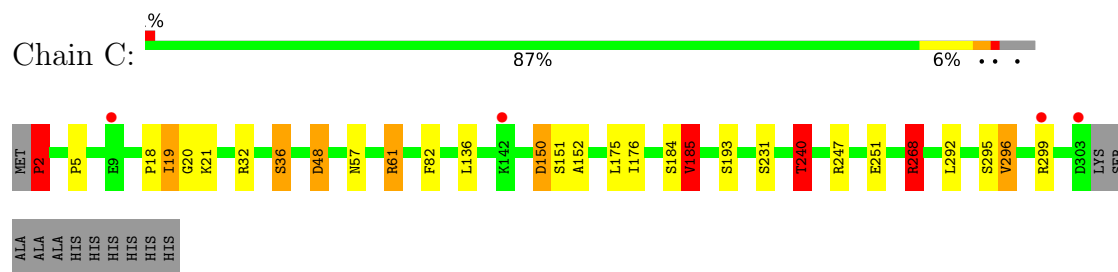
- Molecule 1: Alpha/beta hydrolase fold-3 domain protein



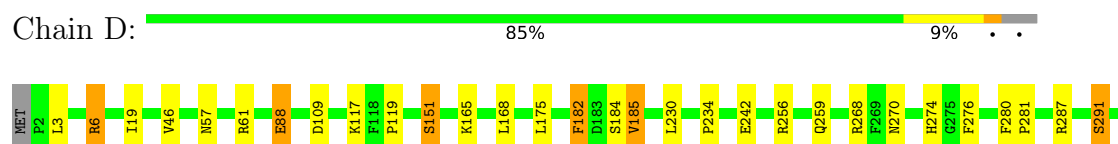
- Molecule 1: Alpha/beta hydrolase fold-3 domain protein

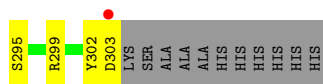


- Molecule 1: Alpha/beta hydrolase fold-3 domain protein



- Molecule 1: Alpha/beta hydrolase fold-3 domain protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	166.41Å 166.41Å 185.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.42 – 2.60 29.42 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (29.42-2.60) 98.3 (29.42-2.60)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.157 , 0.194 0.165 , 0.198	Depositor DCC
R_{free} test set	3952 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9846	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 3.55% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: SO4

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	1.22	6/2397 (0.3%)	1.17	12/3257 (0.4%)
1	B	1.23	5/2397 (0.2%)	1.19	7/3257 (0.2%)
1	C	1.25	6/2397 (0.3%)	1.21	10/3257 (0.3%)
1	D	1.27	10/2397 (0.4%)	1.22	15/3257 (0.5%)
All	All	1.24	27/9588 (0.3%)	1.20	44/13028 (0.3%)

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	D	0	1
All	All	0	2

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	151	SER	CA-C	11.91	1.63	1.53
1	D	151	SER	N-CA	10.08	1.59	1.47
1	C	151	SER	CA-C	9.28	1.61	1.53
1	D	151	SER	CA-CB	9.26	1.67	1.53
1	C	151	SER	CA-CB	8.71	1.69	1.52
1	A	151	SER	CA-CB	8.59	1.69	1.52
1	C	151	SER	N-CA	8.26	1.60	1.47
1	B	151	SER	CB-OG	8.21	1.58	1.42
1	A	151	SER	N-CA	8.04	1.60	1.47
1	C	151	SER	CB-OG	7.53	1.57	1.42
1	A	151	SER	CB-OG	7.45	1.57	1.42
1	A	151	SER	CA-C	7.39	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	88	GLU	CA-CB	7.21	1.64	1.53
1	B	151	SER	N-CA	7.18	1.59	1.47
1	D	151	SER	CB-OG	6.92	1.55	1.42
1	D	151	SER	CA-C	6.46	1.58	1.52
1	B	151	SER	CA-CB	6.27	1.65	1.52
1	C	48	ASP	C-O	-5.94	1.16	1.24
1	D	46	VAL	C-O	-5.89	1.17	1.24
1	D	57	ASN	C-O	-5.70	1.17	1.24
1	C	18	PRO	CA-C	5.66	1.58	1.53
1	D	88	GLU	CG-CD	5.35	1.65	1.52
1	A	234	PRO	CA-C	5.26	1.54	1.51
1	D	165	LYS	CA-C	5.20	1.59	1.52
1	D	234	PRO	CA-C	5.14	1.54	1.51
1	B	228	GLN	C-O	-5.07	1.18	1.24
1	A	151	SER	C-O	5.03	1.29	1.24

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	151	SER	CA-CB-OG	8.44	127.97	111.10
1	D	88	GLU	CB-CG-CD	7.84	125.93	112.60
1	A	151	SER	CA-CB-OG	7.71	126.52	111.10
1	D	88	GLU	CB-CA-C	7.65	123.86	110.85
1	A	151	SER	N-CA-C	-7.25	95.82	108.08
1	B	151	SER	CA-CB-OG	7.18	125.47	111.10
1	C	151	SER	CA-CB-OG	7.08	125.26	111.10
1	A	185	VAL	CB-CA-C	-6.58	104.45	110.63
1	A	234	PRO	N-CA-C	6.29	116.50	110.47
1	B	29	LYS	N-CA-C	-6.23	104.57	111.36
1	D	151	SER	N-CA-C	-6.17	98.37	108.67
1	B	150	ASP	CA-C-O	-6.09	114.23	121.11
1	D	291	SER	N-CA-C	6.08	118.40	111.11
1	B	256	ARG	CB-CA-C	5.92	120.92	110.85
1	C	231	SER	N-CA-C	5.90	117.92	110.24
1	A	185	VAL	N-CA-C	5.86	119.28	113.71
1	D	182	PHE	CB-CA-C	5.82	119.25	109.48
1	D	259	GLN	CA-CB-CG	5.77	125.65	114.10
1	C	295	SER	CB-CA-C	-5.76	101.60	110.81
1	B	193	SER	N-CA-C	5.75	118.28	111.33
1	C	231	SER	CB-CA-C	-5.67	99.60	109.64
1	A	270	ASN	N-CA-C	5.64	118.03	110.35
1	D	185	VAL	N-CA-CB	-5.62	101.95	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	88	GLU	N-CA-C	-5.62	105.24	111.36
1	D	276	PHE	N-CA-C	5.61	118.16	111.71
1	A	72	GLY	N-CA-C	-5.57	104.17	113.02
1	D	19	ILE	CA-C-N	-5.57	110.50	121.41
1	D	19	ILE	C-N-CA	-5.57	110.50	121.41
1	A	150	ASP	CA-C-O	-5.54	114.91	121.16
1	C	150	ASP	CA-C-O	-5.54	114.86	121.11
1	D	268	ARG	CB-CG-CD	-5.53	98.59	111.30
1	D	234	PRO	N-CA-C	5.48	115.73	110.47
1	D	19	ILE	O-C-N	-5.48	116.30	121.94
1	B	22	ALA	N-CA-C	5.46	116.95	110.19
1	C	268	ARG	NE-CZ-NH2	-5.42	114.32	119.20
1	C	185	VAL	N-CA-CB	-5.40	102.33	111.23
1	C	240	THR	N-CA-CB	-5.34	102.39	111.69
1	C	2	PRO	CA-N-CD	-5.21	104.71	112.00
1	B	185	VAL	N-CA-CB	-5.15	102.73	111.23
1	A	58	ILE	N-CA-C	5.08	115.65	107.73
1	A	145	VAL	CB-CA-C	5.06	117.72	110.33
1	A	187	ARG	CG-CD-NE	-5.06	100.88	112.00
1	A	240	THR	CB-CA-C	5.05	120.77	109.56
1	C	193	SER	N-CA-C	5.04	118.20	111.75

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	20	GLY	Peptide
1	D	184	SER	Peptide

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	2342	0	2342	6	0
1	B	2342	0	2342	11	0
1	C	2342	0	2342	18	0
1	D	2342	0	2342	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	15	0	0	0	0
2	D	5	0	0	0	0
3	A	104	0	0	0	0
3	B	118	0	0	0	0
3	C	109	0	0	7	0
3	D	107	0	0	5	0
All	All	9846	0	9368	42	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 2.

All (42) 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:20:GLY:N	3:C:501:HOH:O	1.74	1.14
1:C:36:SER:HB2	3:C:532:HOH:O	1.69	0.91
1:C:19:ILE:O	3:C:502:HOH:O	1.89	0.88
1:C:240:THR:HG21	1:C:251:GLU:OE2	1.82	0.80
1:B:302:TYR:O	1:B:303:ASP:HB2	1.87	0.72
1:D:88:GLU:HG3	3:D:517:HOH:O	1.91	0.68
1:C:19:ILE:CA	3:C:501:HOH:O	2.44	0.66
1:C:48:ASP:OD1	1:C:61:ARG:HG2	1.97	0.65
1:B:302:TYR:O	1:B:303:ASP:CB	2.47	0.62
1:D:256:ARG:HD2	3:D:568:HOH:O	2.00	0.61
1:C:2:PRO:HB3	3:C:590:HOH:O	2.02	0.60
1:A:240:THR:HG21	1:A:251:GLU:OE2	2.01	0.59
1:D:117:LYS:HE3	3:D:558:HOH:O	2.06	0.55
1:D:302:TYR:O	1:D:303:ASP:CG	2.52	0.52
1:C:292:LEU:O	1:C:296:VAL:HG13	2.09	0.52
1:D:117:LYS:CE	3:D:558:HOH:O	2.58	0.51
1:C:240:THR:HG22	1:C:268:ARG:HA	1.91	0.50
1:C:32:ARG:HD3	3:C:602:HOH:O	2.13	0.48
1:B:150:ASP:HA	1:B:176:ILE:O	2.13	0.48
1:D:151:SER:HB2	1:D:274:HIS:NE2	2.28	0.48
1:A:191:GLU:OE2	1:C:2:PRO:HB2	2.14	0.48
1:B:151:SER:HB2	1:B:274:HIS:NE2	2.29	0.47
1:D:230:LEU:CD1	1:D:256:ARG:HG2	2.45	0.47
1:C:251:GLU:OE2	1:C:268:ARG:HD3	2.14	0.46
1:A:267:VAL:HG13	1:B:267:VAL:HG13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ASP:HA	1:A:176:ILE:O	2.16	0.45
1:B:19:ILE:O	1:B:20:GLY:C	2.59	0.45
1:D:242:GLU:HB2	1:D:270:ASN:HA	1.98	0.45
1:A:240:THR:HG22	1:A:268:ARG:HA	1.98	0.45
1:B:142:LYS:HB3	1:B:143:MET:HE3	1.99	0.45
1:C:299:ARG:NH1	1:D:287:ARG:O	2.50	0.45
1:C:150:ASP:HA	1:C:176:ILE:O	2.17	0.44
1:D:6:ARG:NH2	3:D:502:HOH:O	2.39	0.43
1:B:19:ILE:HB	1:B:202:HIS:ND1	2.33	0.43
1:C:82:PHE:CE1	1:C:152:ALA:HB1	2.54	0.42
1:B:91:ASP:HB3	1:B:92:PRO:HD3	2.00	0.42
1:C:32:ARG:NH2	3:C:510:HOH:O	2.53	0.42
1:D:280:PHE:N	1:D:281:PRO:CD	2.83	0.41
1:B:180:VAL:HG22	1:B:249:GLN:HG3	2.03	0.41
1:C:247:ARG:HD2	1:C:268:ARG:HD2	2.01	0.41
1:A:296:VAL:O	1:A:300:THR:HG23	2.20	0.41
1:B:183:ASP:CG	1:C:185:VAL:HG11	2.45	0.41

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	300/314 (96%)	284 (95%)	15 (5%)	1 (0%)	36	58
1	B	300/314 (96%)	283 (94%)	16 (5%)	1 (0%)	36	58
1	C	300/314 (96%)	282 (94%)	17 (6%)	1 (0%)	36	58
1	D	300/314 (96%)	287 (96%)	13 (4%)	0	100	100
All	All	1200/1256 (96%)	1136 (95%)	61 (5%)	3 (0%)	36	58

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ALA
1	B	13	SER
1	C	19	ILE

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	250/259 (96%)	234 (94%)	16 (6%)	16	35
1	B	250/259 (96%)	231 (92%)	19 (8%)	12	27
1	C	250/259 (96%)	237 (95%)	13 (5%)	21	44
1	D	250/259 (96%)	238 (95%)	12 (5%)	23	47
All	All	1000/1036 (96%)	940 (94%)	60 (6%)	17	37

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	5	PRO
1	A	6	ARG
1	A	9	GLU
1	A	29	LYS
1	A	57	ASN
1	A	109	ASP
1	A	115	GLU
1	A	138	LYS
1	A	142	LYS
1	A	145	VAL
1	A	175	LEU
1	A	184	SER
1	A	240	THR
1	A	282	LEU
1	A	300	THR
1	B	3	LEU
1	B	9	GLU
1	B	13	SER

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Mol	Chain	Res	Type
1	B	29	LYS
1	B	34	LEU
1	B	42	GLU
1	B	47	GLU
1	B	87	VAL
1	B	89	SER
1	B	109	ASP
1	B	142	LYS
1	B	143	MET
1	B	175	LEU
1	B	182	PHE
1	B	185	VAL
1	B	187	ARG
1	B	256	ARG
1	B	267	VAL
1	B	303	ASP
1	C	2	PRO
1	C	5	PRO
1	C	21	LYS
1	C	36	SER
1	C	57	ASN
1	C	61	ARG
1	C	136	LEU
1	C	175	LEU
1	C	184	SER
1	C	185	VAL
1	C	240	THR
1	C	268	ARG
1	C	296	VAL
1	D	3	LEU
1	D	6	ARG
1	D	61	ARG
1	D	109	ASP
1	D	119	PRO
1	D	168	LEU
1	D	175	LEU
1	D	182	PHE
1	D	185	VAL
1	D	291	SER
1	D	295	SER
1	D	299	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	ASN
1	A	255	ASN
1	A	271	ASN
1	B	169	ASN
1	B	255	ASN
1	B	259	GLN
1	B	271	ASN
1	C	57	ASN
1	D	134	ASN
1	D	173	GLN
1	D	228	GLN
1	D	259	GLN
1	D	271	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	SO4	B	402	-	4,4,4	0.58	0	6,6,6	0.98	0
2	SO4	C	401	-	4,4,4	0.52	0	6,6,6	0.81	0
2	SO4	A	401	-	4,4,4	0.76	0	6,6,6	1.19	1 (16%)
2	SO4	C	403	-	4,4,4	0.51	0	6,6,6	0.87	0
2	SO4	D	401	-	4,4,4	0.23	0	6,6,6	0.78	0
2	SO4	A	402	-	4,4,4	0.82	0	6,6,6	0.91	0
2	SO4	C	402	-	4,4,4	0.43	0	6,6,6	0.71	0
2	SO4	B	401	-	4,4,4	0.58	0	6,6,6	1.64	1 (16%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	SO4	O4-S-O2	2.53	122.81	109.56
2	A	401	SO4	O3-S-O2	2.06	120.32	109.56

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 [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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	302/314 (96%)	-0.63	2 (0%) 84 82	15, 26, 49, 65	0
1	B	302/314 (96%)	-0.68	4 (1%) 75 71	16, 24, 50, 72	0
1	C	302/314 (96%)	-0.57	4 (1%) 75 71	16, 26, 46, 73	0
1	D	302/314 (96%)	-0.73	1 (0%) 90 88	15, 23, 45, 69	0
All	All	1208/1256 (96%)	-0.65	11 (0%) 81 78	15, 24, 47, 73	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	299	ARG	4.7
1	D	303	ASP	4.3
1	B	303	ASP	3.9
1	C	303	ASP	3.7
1	B	21	LYS	3.2
1	B	142	LYS	3.0
1	B	36	SER	2.9
1	A	256	ARG	2.1
1	A	303	ASP	2.1
1	C	9	GLU	2.1
1	C	142	LYS	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	SO4	A	402	5/5	0.95	0.13	53,53,55,60	0
2	SO4	B	402	5/5	0.95	0.10	55,56,60,65	0
2	SO4	C	403	5/5	0.95	0.12	42,49,52,57	0
2	SO4	C	401	5/5	0.96	0.11	42,45,54,57	0
2	SO4	B	401	5/5	0.99	0.07	34,34,36,36	0
2	SO4	C	402	5/5	0.99	0.04	31,31,34,35	0
2	SO4	A	401	5/5	0.99	0.06	31,32,33,37	0
2	SO4	D	401	5/5	0.99	0.05	28,29,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.