



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2024 – 06:00 PM EST

PDB ID : 2QUB
Title : Crystal structure of extracellular lipase LipA from *Serratia marcescens*
Authors : Meier, R.; Baumann, U.
Deposited on : 2007-08-04
Resolution : 1.80 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

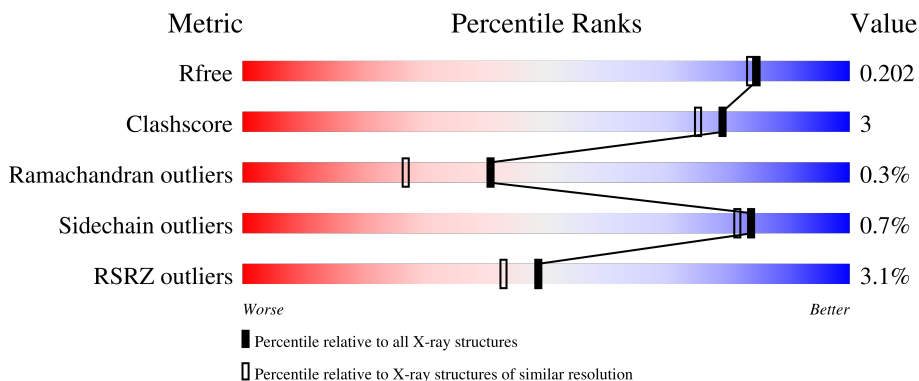
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



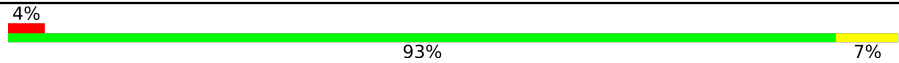
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	615	 3% 96%
1	C	615	 2% 95%
1	E	615	 3% 93% 7%
1	G	615	 3% 95% 5%
1	I	615	 3% 95% 5%

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Mol	Chain	Length	Quality of chain
1	K	615	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a small red segment on the left labeled '4%', a large green segment in the middle labeled '93%', and a small yellow segment on the right labeled '7%'.</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 29575 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 Extracellular lipase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	615	4601	2883	779	936	3	0	0	0
1	C	615	4601	2883	779	936	3	0	0	0
1	E	615	4601	2883	779	936	3	0	0	0
1	G	615	4601	2883	779	936	3	0	0	0
1	I	615	4601	2883	779	936	3	0	0	0
1	K	615	4601	2883	779	936	3	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP Q59933
A	0	HIS	-	expression tag	UNP Q59933
C	-1	SER	-	expression tag	UNP Q59933
C	0	HIS	-	expression tag	UNP Q59933
E	-1	SER	-	expression tag	UNP Q59933
E	0	HIS	-	expression tag	UNP Q59933
G	-1	SER	-	expression tag	UNP Q59933
G	0	HIS	-	expression tag	UNP Q59933
I	-1	SER	-	expression tag	UNP Q59933
I	0	HIS	-	expression tag	UNP Q59933
K	-1	SER	-	expression tag	UNP Q59933
K	0	HIS	-	expression tag	UNP Q59933

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total Ca 8 8	0	0
2	C	8	Total Ca 8 8	0	0
2	E	8	Total Ca 8 8	0	0
2	G	8	Total Ca 8 8	0	0
2	I	8	Total Ca 8 8	0	0
2	K	8	Total Ca 8 8	0	0

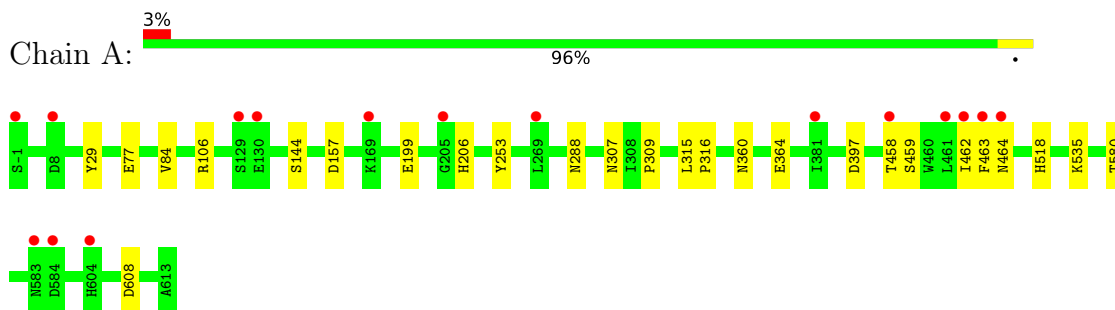
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	322	Total O 322 322	0	0
3	C	360	Total O 360 360	0	0
3	E	285	Total O 285 285	0	0
3	G	346	Total O 346 346	0	0
3	I	305	Total O 305 305	0	0
3	K	303	Total O 303 303	0	0

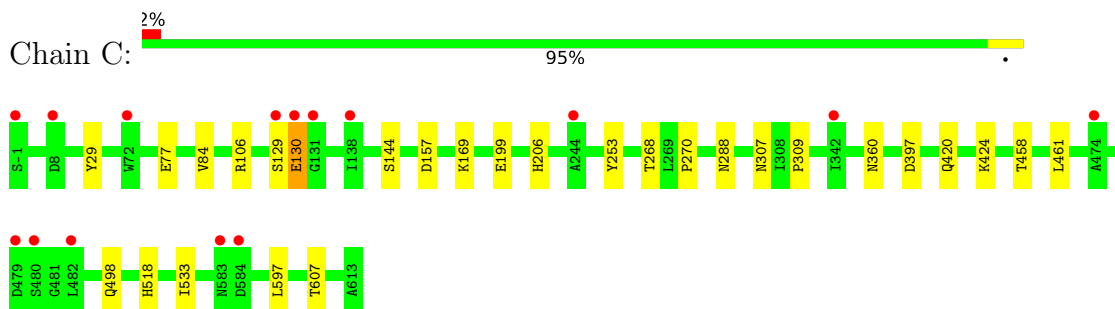
3 Residue-property plots [i](#)

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.

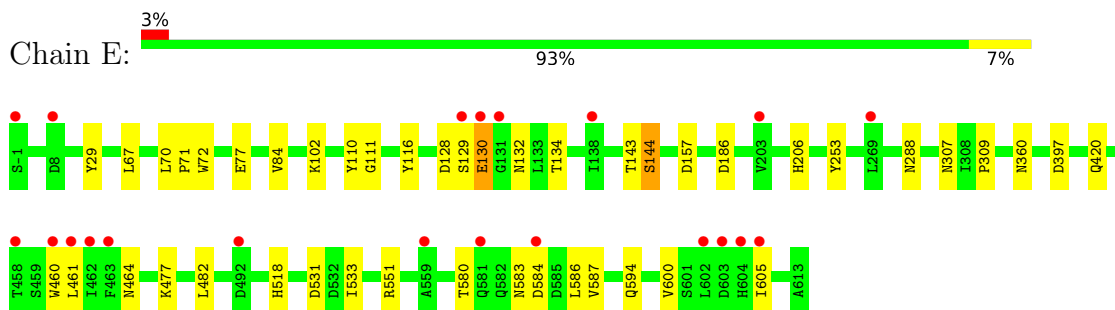
- Molecule 1: Extracellular lipase



- Molecule 1: Extracellular lipase

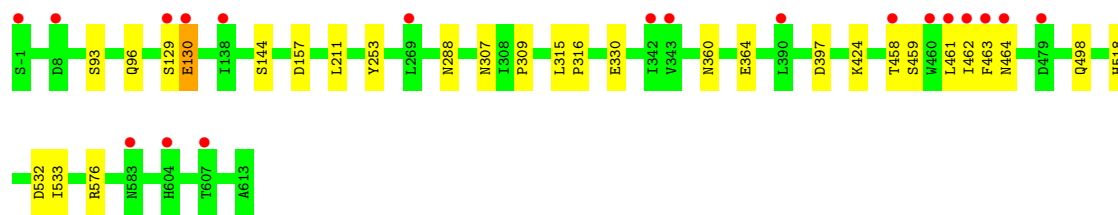


- Molecule 1: Extracellular lipase

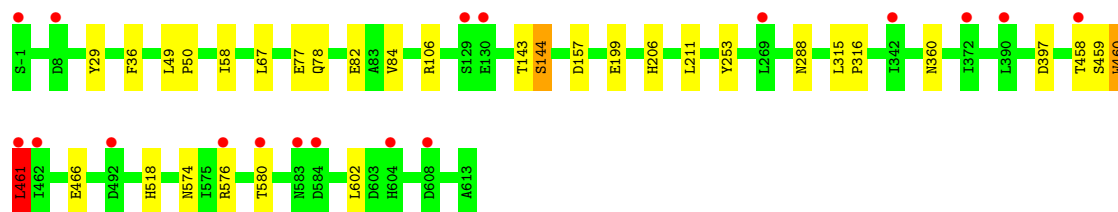
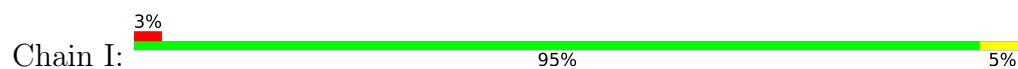


- Molecule 1: Extracellular lipase

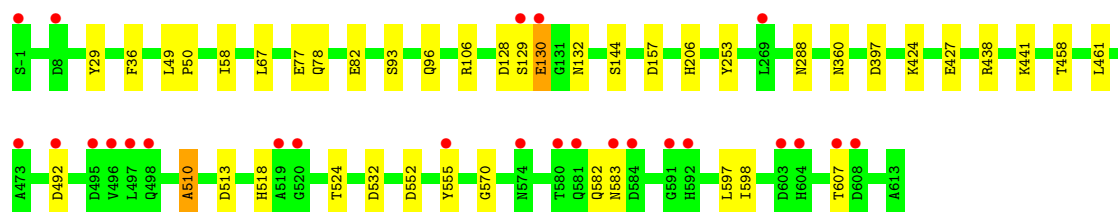
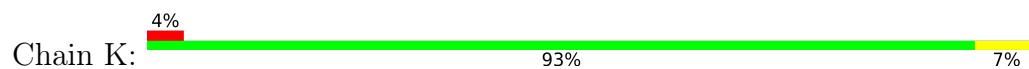




- Molecule 1: Extracellular lipase



- Molecule 1: Extracellular lipase



4 Data and refinement statistics i

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	202.40Å 202.40Å 317.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.93 – 1.80 24.83 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.6 (24.93-1.80) 95.6 (24.83-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 1.80Å)	Xtrriage
Refinement program	REFMAC 5.3.0036	Depositor
R, R_{free}	0.171 , 0.200 0.174 , 0.202	Depositor DCC
R_{free} test set	1275 reflections (0.30%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.015 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	29575	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.53 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.8798e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.46	0/4699	0.56	2/6387 (0.0%)
1	C	0.50	0/4699	0.57	2/6387 (0.0%)
1	E	0.45	0/4699	0.56	2/6387 (0.0%)
1	G	0.48	0/4699	0.58	3/6387 (0.0%)
1	I	0.48	0/4699	0.57	3/6387 (0.0%)
1	K	0.47	0/4699	0.57	3/6387 (0.0%)
All	All	0.47	0/28194	0.57	15/38322 (0.0%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	157	ASP	CB-CG-OD1	7.87	125.38	118.30
1	K	157	ASP	CB-CG-OD1	6.96	124.56	118.30
1	G	157	ASP	CB-CG-OD1	6.85	124.46	118.30
1	E	397	ASP	CB-CG-OD2	6.74	124.37	118.30
1	K	397	ASP	CB-CG-OD2	6.62	124.25	118.30
1	E	157	ASP	CB-CG-OD1	6.62	124.25	118.30
1	A	157	ASP	CB-CG-OD1	6.48	124.14	118.30
1	C	157	ASP	CB-CG-OD1	6.33	124.00	118.30
1	C	397	ASP	CB-CG-OD2	5.58	123.32	118.30
1	G	397	ASP	CB-CG-OD2	5.48	123.23	118.30
1	I	461	LEU	N-CA-C	5.46	125.73	111.00
1	G	532	ASP	CB-CG-OD2	5.42	123.18	118.30
1	A	397	ASP	CB-CG-OD2	5.19	122.97	118.30
1	K	552	ASP	CB-CG-OD1	5.15	122.94	118.30
1	I	397	ASP	CB-CG-OD2	5.12	122.91	118.30

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	4601	0	4339	25	0
1	C	4601	0	4339	25	2
1	E	4601	0	4339	30	0
1	G	4601	0	4339	26	0
1	I	4601	0	4339	21	0
1	K	4601	0	4339	27	0
2	A	8	0	0	0	0
2	C	8	0	0	0	0
2	E	8	0	0	0	0
2	G	8	0	0	0	0
2	I	8	0	0	0	0
2	K	8	0	0	0	0
3	A	322	0	0	10	0
3	C	360	0	0	13	0
3	E	285	0	0	8	0
3	G	346	0	0	7	0
3	I	305	0	0	4	0
3	K	303	0	0	6	2
All	All	29575	0	26034	152	2

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 3.

All (152) 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:199:GLU:HG3	3:C:923:HOH:O	1.52	1.10
1:K:424:LYS:HE3	3:K:763:HOH:O	1.53	1.05
1:E:129:SER:O	1:E:130:GLU:HG3	1.59	1.02
1:I:461:LEU:O	1:I:461:LEU:HG	1.53	1.01
1:A:458:THR:HG23	1:A:463:PHE:HA	1.46	0.98
1:C:424:LYS:CE	3:C:839:HOH:O	2.16	0.93
1:I:458:THR:HG23	1:I:461:LEU:HD12	1.54	0.89
1:K:129:SER:O	1:K:130:GLU:HG3	1.73	0.88
1:C:129:SER:O	1:C:130:GLU:HG3	1.71	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:424:LYS:HE3	3:C:867:HOH:O	1.72	0.88
1:I:458:THR:CG2	1:I:461:LEU:HD12	2.08	0.83
1:E:600:VAL:HG11	1:E:605:ILE:HD11	1.63	0.81
1:K:438:ARG:NH2	1:K:524:THR:OG1	2.15	0.80
1:A:535:LYS:HE2	3:A:874:HOH:O	1.82	0.79
1:G:129:SER:O	1:G:130:GLU:HG3	1.83	0.78
1:K:427:GLU:OE2	1:K:438:ARG:HD2	1.85	0.77
1:G:364:GLU:OE2	3:G:924:HOH:O	2.03	0.76
1:G:458:THR:HG23	1:G:463:PHE:HA	1.67	0.76
1:K:424:LYS:HD2	3:K:825:HOH:O	1.86	0.76
1:K:458:THR:HG21	1:K:461:LEU:O	1.87	0.75
1:A:458:THR:CG2	1:A:463:PHE:HA	2.18	0.73
1:K:424:LYS:CE	3:K:825:HOH:O	2.36	0.73
1:E:307:ASN:OD1	1:E:309:PRO:HD2	1.89	0.73
1:A:463:PHE:O	1:A:464:ASN:ND2	2.22	0.72
1:C:424:LYS:CD	3:C:839:HOH:O	2.33	0.72
1:C:169:LYS:HE2	3:C:747:HOH:O	1.88	0.72
1:K:424:LYS:CD	3:K:825:HOH:O	2.38	0.72
1:C:424:LYS:HD2	3:C:839:HOH:O	1.89	0.71
1:C:169:LYS:NZ	3:C:747:HOH:O	2.23	0.70
1:K:129:SER:O	1:K:130:GLU:CG	2.39	0.70
1:G:129:SER:O	1:G:130:GLU:CG	2.38	0.70
1:I:576:ARG:HG2	1:I:602:LEU:HD21	1.73	0.70
1:E:587:VAL:HG23	3:E:888:HOH:O	1.92	0.69
1:E:586:LEU:C	3:E:888:HOH:O	2.30	0.69
1:G:463:PHE:O	1:G:464:ASN:ND2	2.26	0.69
1:G:307:ASN:OD1	1:G:309:PRO:HD2	1.92	0.68
1:G:129:SER:O	1:G:130:GLU:CB	2.39	0.68
1:I:199:GLU:HG3	3:I:792:HOH:O	1.94	0.68
1:A:458:THR:HG23	1:A:463:PHE:CA	2.22	0.68
1:A:580:THR:HG23	3:A:862:HOH:O	1.93	0.68
1:C:420:GLN:OE1	3:C:899:HOH:O	2.12	0.67
1:C:77:GLU:OE1	1:C:106:ARG:CZ	2.42	0.67
1:C:169:LYS:CE	3:C:747:HOH:O	2.42	0.67
1:C:458:THR:HG21	1:C:461:LEU:O	1.95	0.66
1:K:607:THR:HG23	3:K:852:HOH:O	1.96	0.66
1:E:71:PRO:HD2	1:E:72:TRP:CZ3	2.31	0.65
1:E:77:GLU:HG3	3:E:809:HOH:O	1.95	0.65
1:E:420:GLN:OE1	3:E:853:HOH:O	2.14	0.65
1:A:458:THR:HG22	1:A:459:SER:N	2.12	0.64
1:C:129:SER:O	1:C:130:GLU:CG	2.44	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:GLU:OE1	1:A:106:ARG:CZ	2.45	0.64
1:C:498:GLN:HG3	3:C:786:HOH:O	1.96	0.64
1:G:498:GLN:CG	3:G:919:HOH:O	2.45	0.64
1:E:586:LEU:CD2	1:E:605:ILE:HG13	2.28	0.62
1:K:129:SER:O	1:K:130:GLU:CB	2.48	0.61
1:E:29:TYR:OH	1:E:206:HIS:HD2	1.83	0.61
1:G:498:GLN:HG3	3:G:919:HOH:O	2.01	0.61
1:G:458:THR:HG22	1:G:459:SER:N	2.17	0.60
1:K:78:GLN:NE2	1:K:82:GLU:OE2	2.31	0.60
1:A:459:SER:N	3:A:907:HOH:O	2.33	0.60
1:K:93:SER:H	1:K:96:GLN:HE21	1.49	0.60
1:A:458:THR:HG22	1:A:459:SER:H	1.66	0.60
1:C:307:ASN:OD1	1:C:309:PRO:HD2	2.03	0.59
1:I:574:ASN:HD21	1:I:576:ARG:NH1	2.00	0.59
1:K:58:ILE:HG22	1:K:67:LEU:HD22	1.84	0.59
1:I:78:GLN:NE2	1:I:82:GLU:OE2	2.34	0.58
1:A:307:ASN:OD1	1:A:309:PRO:HD2	2.02	0.58
1:K:77:GLU:OE1	1:K:106:ARG:CZ	2.50	0.58
1:A:199:GLU:HG3	3:A:872:HOH:O	2.02	0.58
1:G:129:SER:O	1:G:130:GLU:HB2	2.03	0.58
1:I:466:GLU:HG3	3:I:912:HOH:O	2.02	0.58
1:C:498:GLN:CG	3:C:786:HOH:O	2.51	0.58
1:I:77:GLU:OE1	1:I:106:ARG:CZ	2.52	0.57
1:A:364:GLU:OE2	3:A:942:HOH:O	2.18	0.56
1:I:574:ASN:HD21	1:I:576:ARG:HH11	1.52	0.56
1:K:29:TYR:OH	1:K:206:HIS:HD2	1.89	0.56
1:A:29:TYR:OH	1:A:206:HIS:HD2	1.88	0.56
1:E:580:THR:HG23	3:E:832:HOH:O	2.06	0.56
1:A:462:ILE:HG13	1:A:462:ILE:O	2.06	0.55
1:K:582:GLN:O	1:K:583:ASN:HB2	2.08	0.54
1:C:424:LYS:HE2	3:C:839:HOH:O	1.93	0.54
1:G:458:THR:HG23	1:G:463:PHE:CA	2.35	0.54
1:I:580:THR:HG23	3:I:881:HOH:O	2.07	0.54
1:C:29:TYR:OH	1:C:206:HIS:HD2	1.90	0.54
1:I:58:ILE:HG22	1:I:67:LEU:HD22	1.89	0.54
1:A:458:THR:C	3:A:907:HOH:O	2.46	0.53
1:I:29:TYR:OH	1:I:206:HIS:HD2	1.93	0.52
1:G:462:ILE:O	1:G:462:ILE:HG13	2.10	0.51
1:G:458:THR:CG2	1:G:463:PHE:HA	2.38	0.51
1:G:498:GLN:HG2	3:G:919:HOH:O	2.05	0.51
1:E:533:ILE:HD12	1:E:533:ILE:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:459:SER:CB	3:A:907:HOH:O	2.59	0.50
1:G:458:THR:HG23	1:G:463:PHE:C	2.31	0.50
1:E:128:ASP:OD2	1:E:132:ASN:HB2	2.12	0.49
1:A:535:LYS:CE	3:A:874:HOH:O	2.53	0.49
1:I:461:LEU:O	1:I:461:LEU:CG	2.37	0.49
1:C:607:THR:HG21	3:I:873:HOH:O	2.12	0.48
1:G:93:SER:OG	1:G:96:GLN:HG3	2.14	0.48
1:E:600:VAL:HG11	1:E:605:ILE:CD1	2.39	0.48
1:K:555:TYR:CZ	1:K:598:ILE:HD12	2.49	0.48
1:E:129:SER:O	1:E:130:GLU:CG	2.48	0.47
1:I:458:THR:HG21	1:I:461:LEU:HD12	1.92	0.47
1:E:587:VAL:N	3:E:888:HOH:O	2.46	0.47
1:G:461:LEU:HD22	3:G:903:HOH:O	2.14	0.47
1:C:268:THR:OG1	1:C:270:PRO:HD2	2.14	0.47
1:E:186:ASP:OD1	1:G:576:ARG:NH2	2.48	0.47
1:A:458:THR:HG23	1:A:463:PHE:C	2.34	0.46
1:G:424:LYS:HD2	3:G:868:HOH:O	2.15	0.46
1:K:49:LEU:N	1:K:50:PRO:HD2	2.31	0.46
1:A:459:SER:HB3	3:A:907:HOH:O	2.17	0.45
1:A:458:THR:CG2	1:A:459:SER:N	2.80	0.45
1:E:460:TRP:O	1:E:461:LEU:HB2	2.17	0.45
1:I:143:THR:O	1:I:144:SER:CB	2.65	0.44
1:E:143:THR:O	1:E:144:SER:CB	2.65	0.44
1:I:211:LEU:C	1:I:211:LEU:HD23	2.37	0.44
1:I:459:SER:O	1:I:460:TRP:C	2.54	0.44
1:A:608:ASP:OD1	1:A:608:ASP:N	2.51	0.44
1:G:211:LEU:HD23	1:G:211:LEU:C	2.37	0.44
1:E:531:ASP:HB3	1:E:551:ARG:O	2.18	0.44
1:C:129:SER:O	1:C:130:GLU:CB	2.65	0.44
1:E:102:LYS:HE3	1:E:110:TYR:CD1	2.53	0.44
1:E:111:GLY:HA3	1:E:116:TYR:O	2.18	0.43
1:K:513:ASP:HB2	1:K:532:ASP:OD1	2.18	0.43
1:A:253:TYR:CG	1:A:288:ASN:HB2	2.53	0.43
1:C:597:LEU:N	1:C:597:LEU:HD12	2.33	0.43
1:E:67:LEU:HD23	1:E:70:LEU:HD11	2.01	0.43
1:G:330:GLU:HG3	3:G:828:HOH:O	2.19	0.43
1:G:253:TYR:CG	1:G:288:ASN:HB2	2.54	0.43
1:I:253:TYR:CG	1:I:288:ASN:HB2	2.53	0.43
1:C:253:TYR:CG	1:C:288:ASN:HB2	2.54	0.42
1:K:128:ASP:OD2	1:K:132:ASN:HB2	2.19	0.42
1:K:441:LYS:HB3	1:K:441:LYS:HE2	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:492:ASP:HA	1:K:510:ALA:HB3	2.01	0.42
1:E:464:ASN:HD22	1:E:464:ASN:N	2.18	0.42
1:K:129:SER:O	1:K:130:GLU:HB2	2.19	0.42
1:E:477:LYS:HG2	1:E:482:LEU:HD13	2.01	0.42
1:C:533:ILE:HD11	1:G:533:ILE:HD11	2.01	0.42
1:E:586:LEU:N	3:E:888:HOH:O	2.51	0.42
1:E:253:TYR:CG	1:E:288:ASN:HB2	2.55	0.42
1:C:424:LYS:NZ	3:C:839:HOH:O	2.35	0.42
1:K:253:TYR:CG	1:K:288:ASN:HB2	2.55	0.42
1:G:458:THR:CG2	1:G:459:SER:N	2.83	0.41
1:K:597:LEU:N	1:K:597:LEU:HD12	2.35	0.41
1:A:462:ILE:HG13	3:A:907:HOH:O	2.20	0.41
1:A:315:LEU:HA	1:A:316:PRO:HD3	1.99	0.41
1:I:315:LEU:HA	1:I:316:PRO:HD3	1.96	0.41
1:E:551:ARG:HG2	1:E:594:GLN:HG2	2.03	0.40
1:E:583:ASN:HB3	1:E:584:ASP:H	1.65	0.40
1:K:570:GLY:O	3:K:820:HOH:O	2.21	0.40
1:E:587:VAL:CG2	3:E:888:HOH:O	2.61	0.40
1:I:49:LEU:N	1:I:50:PRO:CD	2.85	0.40
1:G:315:LEU:HA	1:G:316:PRO:HD3	1.92	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:461:LEU:CD1	3:K:768:HOH:O[9_444]	1.05	1.15
1:C:461:LEU:CG	3:K:768:HOH:O[9_444]	2.07	0.13

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	613/615 (100%)	602 (98%)	10 (2%)	1 (0%)	47	33
1	C	613/615 (100%)	599 (98%)	12 (2%)	2 (0%)	41	27
1	E	613/615 (100%)	598 (98%)	13 (2%)	2 (0%)	41	27
1	G	613/615 (100%)	600 (98%)	11 (2%)	2 (0%)	41	27
1	I	613/615 (100%)	601 (98%)	10 (2%)	2 (0%)	41	27
1	K	613/615 (100%)	592 (97%)	18 (3%)	3 (0%)	29	15
All	All	3678/3690 (100%)	3592 (98%)	74 (2%)	12 (0%)	41	27

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	144	SER
1	C	130	GLU
1	C	144	SER
1	E	130	GLU
1	E	144	SER
1	G	130	GLU
1	G	144	SER
1	I	144	SER
1	I	461	LEU
1	K	130	GLU
1	K	144	SER
1	K	510	ALA

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	474/474 (100%)	471 (99%)	3 (1%)	86	84
1	C	474/474 (100%)	471 (99%)	3 (1%)	86	84
1	E	474/474 (100%)	470 (99%)	4 (1%)	81	78
1	G	474/474 (100%)	472 (100%)	2 (0%)	91	89
1	I	474/474 (100%)	469 (99%)	5 (1%)	73	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	K	474/474 (100%)	471 (99%)	3 (1%)	86	84
All	All	2844/2844 (100%)	2824 (99%)	20 (1%)	84	81

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	VAL
1	A	360	ASN
1	A	518	HIS
1	C	84	VAL
1	C	360	ASN
1	C	518	HIS
1	E	84	VAL
1	E	134	THR
1	E	360	ASN
1	E	518	HIS
1	G	360	ASN
1	G	518	HIS
1	I	36	PHE
1	I	84	VAL
1	I	360	ASN
1	I	460	TRP
1	I	518	HIS
1	K	36	PHE
1	K	360	ASN
1	K	518	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	206	HIS
1	A	275	HIS
1	A	328	ASN
1	A	464	ASN
1	A	498	GLN
1	C	182	ASN
1	C	206	HIS
1	C	420	GLN
1	C	433	ASN
1	C	464	ASN

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Mol	Chain	Res	Type
1	C	522	ASN
1	C	582	GLN
1	C	583	ASN
1	E	85	ASN
1	E	206	HIS
1	E	275	HIS
1	E	420	GLN
1	E	464	ASN
1	E	522	ASN
1	G	275	HIS
1	G	362	ASN
1	G	433	ASN
1	G	464	ASN
1	I	206	HIS
1	I	328	ASN
1	I	433	ASN
1	I	498	GLN
1	I	582	GLN
1	K	62	GLN
1	K	96	GLN
1	K	206	HIS
1	K	498	GLN
1	K	522	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 48 ligands modelled in this entry, 48 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

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	615/615 (100%)	-0.03	16 (2%) 56 51	14, 22, 31, 43	0
1	C	615/615 (100%)	-0.07	15 (2%) 59 54	14, 22, 29, 43	0
1	E	615/615 (100%)	0.11	21 (3%) 45 39	13, 22, 32, 41	0
1	G	615/615 (100%)	-0.03	19 (3%) 49 43	13, 22, 30, 42	0
1	I	615/615 (100%)	-0.03	18 (2%) 51 46	16, 22, 30, 44	0
1	K	615/615 (100%)	0.18	25 (4%) 37 31	15, 22, 30, 44	0
All	All	3690/3690 (100%)	0.02	114 (3%) 49 43	13, 22, 30, 44	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	583	ASN	7.4
1	K	584	ASP	6.7
1	E	129	SER	6.5
1	G	463	PHE	6.3
1	E	462	ILE	5.1
1	K	129	SER	4.8
1	I	461	LEU	4.7
1	K	130	GLU	4.5
1	C	129	SER	4.3
1	E	460	TRP	4.2
1	I	269	LEU	4.1
1	I	129	SER	4.0
1	A	583	ASN	3.9
1	K	607	THR	3.9
1	E	8	ASP	3.9
1	G	130	GLU	3.8
1	E	130	GLU	3.8
1	E	463	PHE	3.7
1	K	-1	SER	3.7

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Mol	Chain	Res	Type	RSRZ
1	K	492	ASP	3.7
1	G	461	LEU	3.6
1	A	464	ASN	3.6
1	K	603	ASP	3.5
1	K	496	VAL	3.4
1	C	130	GLU	3.4
1	A	130	GLU	3.4
1	E	203	VAL	3.2
1	K	473	ALA	3.2
1	I	576	ARG	3.2
1	E	584	ASP	3.1
1	G	129	SER	3.1
1	K	592	HIS	3.1
1	I	604	HIS	3.1
1	A	8	ASP	3.1
1	A	129	SER	3.1
1	K	580	THR	3.1
1	A	463	PHE	3.0
1	A	462	ILE	3.0
1	K	269	LEU	3.0
1	G	460	TRP	2.9
1	E	604	HIS	2.9
1	E	131	GLY	2.9
1	E	603	ASP	2.9
1	C	131	GLY	2.9
1	I	583	ASN	2.8
1	E	269	LEU	2.8
1	C	583	ASN	2.8
1	K	604	HIS	2.8
1	K	581	GLN	2.8
1	G	458	THR	2.7
1	A	461	LEU	2.7
1	I	492	ASP	2.7
1	C	244	ALA	2.7
1	K	520	GLY	2.7
1	K	591	GLY	2.7
1	A	584	ASP	2.6
1	K	574	ASN	2.6
1	E	461	LEU	2.6
1	I	8	ASP	2.6
1	E	559	ALA	2.6
1	I	130	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	605	ILE	2.6
1	C	-1	SER	2.6
1	I	584	ASP	2.6
1	C	479	ASP	2.5
1	G	8	ASP	2.5
1	A	458	THR	2.5
1	G	342	ILE	2.5
1	I	462	ILE	2.5
1	C	8	ASP	2.5
1	E	-1	SER	2.5
1	I	458	THR	2.5
1	A	-1	SER	2.5
1	G	604	HIS	2.4
1	G	479	ASP	2.4
1	K	555	TYR	2.4
1	K	519	ALA	2.4
1	A	269	LEU	2.4
1	K	8	ASP	2.4
1	K	498	GLN	2.4
1	G	583	ASN	2.4
1	C	72	TRP	2.4
1	C	584	ASP	2.4
1	E	138	ILE	2.4
1	E	492	ASP	2.3
1	I	342	ILE	2.3
1	A	604	HIS	2.3
1	G	269	LEU	2.3
1	I	-1	SER	2.2
1	E	458	THR	2.2
1	C	138	ILE	2.1
1	C	482	LEU	2.1
1	K	495	ASP	2.1
1	C	480	SER	2.1
1	G	462	ILE	2.1
1	G	464	ASN	2.1
1	K	497	LEU	2.1
1	C	474	ALA	2.1
1	G	343	VAL	2.1
1	G	390	LEU	2.1
1	A	169	LYS	2.1
1	A	205	GLY	2.1
1	E	581	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	607	THR	2.1
1	K	608	ASP	2.1
1	A	381	ILE	2.1
1	C	342	ILE	2.1
1	I	580	THR	2.0
1	G	138	ILE	2.0
1	I	608	ASP	2.0
1	G	-1	SER	2.0
1	E	602	LEU	2.0
1	I	372	ILE	2.0
1	I	390	LEU	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 monosaccharides 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	CA	E	615	1/1	0.91	0.12	27,27,27,27	0
2	CA	E	616	1/1	0.94	0.10	28,28,28,28	0
2	CA	E	614	1/1	0.96	0.13	28,28,28,28	0
2	CA	K	614	1/1	0.96	0.19	28,28,28,28	0
2	CA	A	614	1/1	0.97	0.09	23,23,23,23	0
2	CA	K	616	1/1	0.97	0.17	25,25,25,25	0
2	CA	I	615	1/1	0.98	0.08	24,24,24,24	0
2	CA	I	616	1/1	0.98	0.12	25,25,25,25	0
2	CA	A	619	1/1	0.98	0.04	16,16,16,16	0
2	CA	K	615	1/1	0.98	0.18	25,25,25,25	0
2	CA	I	614	1/1	0.98	0.12	28,28,28,28	0
2	CA	K	619	1/1	0.98	0.04	17,17,17,17	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CA	E	619	1/1	0.99	0.05	15,15,15,15	0
2	CA	E	620	1/1	0.99	0.07	14,14,14,14	0
2	CA	G	614	1/1	0.99	0.07	20,20,20,20	0
2	CA	G	616	1/1	0.99	0.05	18,18,18,18	0
2	CA	G	619	1/1	0.99	0.03	16,16,16,16	0
2	CA	G	620	1/1	0.99	0.03	13,13,13,13	0
2	CA	G	621	1/1	0.99	0.05	15,15,15,15	0
2	CA	C	614	1/1	0.99	0.08	20,20,20,20	0
2	CA	C	619	1/1	0.99	0.06	15,15,15,15	0
2	CA	C	621	1/1	0.99	0.05	15,15,15,15	0
2	CA	I	619	1/1	0.99	0.03	16,16,16,16	0
2	CA	I	621	1/1	0.99	0.04	14,14,14,14	0
2	CA	A	616	1/1	0.99	0.07	21,21,21,21	0
2	CA	A	615	1/1	0.99	0.06	20,20,20,20	0
2	CA	A	621	1/1	0.99	0.06	17,17,17,17	0
2	CA	K	617	1/1	0.99	0.06	18,18,18,18	0
2	CA	K	618	1/1	0.99	0.06	19,19,19,19	0
2	CA	E	617	1/1	0.99	0.04	15,15,15,15	0
2	CA	K	620	1/1	0.99	0.06	15,15,15,15	0
2	CA	K	621	1/1	0.99	0.06	15,15,15,15	0
2	CA	C	617	1/1	1.00	0.02	14,14,14,14	0
2	CA	E	618	1/1	1.00	0.06	16,16,16,16	0
2	CA	C	618	1/1	1.00	0.04	16,16,16,16	0
2	CA	I	617	1/1	1.00	0.05	16,16,16,16	0
2	CA	I	618	1/1	1.00	0.04	17,17,17,17	0
2	CA	A	620	1/1	1.00	0.04	14,14,14,14	0
2	CA	I	620	1/1	1.00	0.04	15,15,15,15	0
2	CA	E	621	1/1	1.00	0.03	15,15,15,15	0
2	CA	C	620	1/1	1.00	0.04	13,13,13,13	0
2	CA	G	615	1/1	1.00	0.05	17,17,17,17	0
2	CA	A	618	1/1	1.00	0.06	16,16,16,16	0
2	CA	G	617	1/1	1.00	0.04	15,15,15,15	0
2	CA	G	618	1/1	1.00	0.06	17,17,17,17	0
2	CA	A	617	1/1	1.00	0.03	14,14,14,14	0
2	CA	C	615	1/1	1.00	0.06	16,16,16,16	0
2	CA	C	616	1/1	1.00	0.05	18,18,18,18	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.