



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:11 AM UTC

PDB ID : 1L3F / pdb_00001l3f
Title : Thermolysin in the Absence of Substrate has an Open Conformation
Authors : Hausrath, A.C.; Matthews, B.W.
Deposited on : 2002-02-26
Resolution : 2.30 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
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

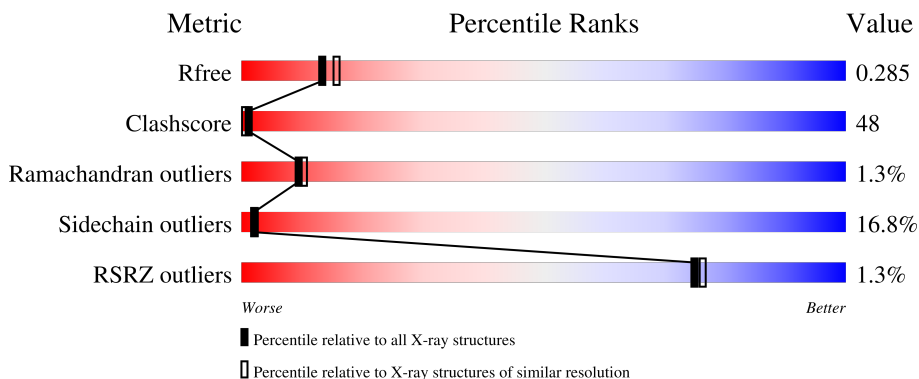
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	E	316	28% 41% 27% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2561 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 Thermolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	316	Total	C	N	O	S	0	0	0
			2427	1525	406	494	2			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	4	Total	Ca	0	0
			4	4		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	3	Total	Zn	0	0
			3	3		

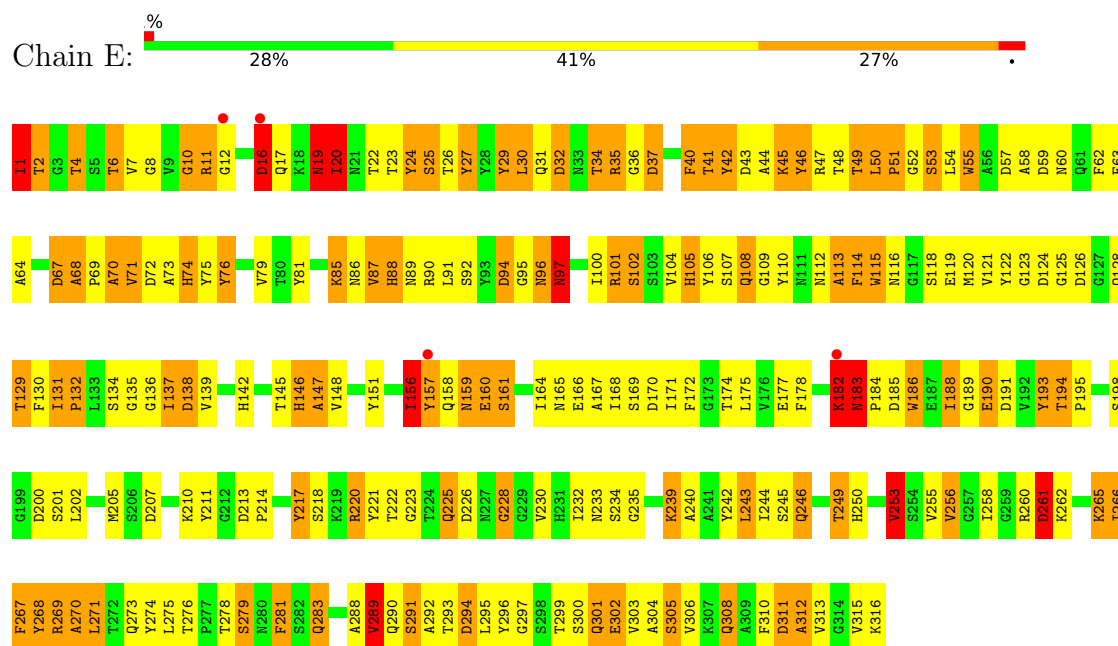
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	127	Total	O	0	0
			127	127		

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: Thermolysin



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	97.05Å 97.05Å 106.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 25.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.7 (25.00-2.30) 96.7 (25.00-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.31Å)	Xtriage
Refinement program		Depositor
R, R_{free}	0.202 , 0.302 0.196 , 0.285	Depositor DCC
R_{free} test set	1138 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 143.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2561	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.41% 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: CA, ZN

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	E	1.47	11/2486 (0.4%)	2.26	132/3386 (3.9%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	74	HIS	N-CA	-6.79	1.38	1.46
1	E	131	ILE	CA-CB	-6.32	1.48	1.55
1	E	16	ASP	CG-OD2	5.93	1.36	1.25
1	E	131	ILE	N-CA	-5.85	1.41	1.46
1	E	226	ASP	CG-OD2	5.48	1.35	1.25
1	E	20	ILE	CA-CB	-5.47	1.49	1.55
1	E	302	GLU	CD-OE2	5.38	1.35	1.25
1	E	311	ASP	CG-OD2	5.10	1.35	1.25
1	E	194	THR	CA-C	-5.05	1.46	1.52
1	E	289	VAL	C-N	-5.02	1.27	1.33
1	E	268	TYR	N-CA	-5.01	1.40	1.46

All (132) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	261	ASP	CA-CB-CG	-9.97	102.63	112.60
1	E	156	ILE	N-CA-CB	-9.88	99.13	110.99
1	E	96	ASN	CA-CB-CG	8.46	121.06	112.60
1	E	301	GLN	CA-C-O	8.30	129.22	120.42
1	E	35	ARG	N-CA-C	8.26	123.19	108.24
1	E	94	ASP	CA-CB-CG	-8.23	104.37	112.60
1	E	138	ASP	CA-CB-CG	8.13	120.73	112.60
1	E	289	VAL	CA-C-N	-7.98	108.96	120.29
1	E	289	VAL	C-N-CA	-7.98	108.96	120.29
1	E	75	TYR	N-CA-C	7.78	119.39	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	94	ASP	N-CA-C	-7.69	103.48	113.17
1	E	1	ILE	CA-C-N	7.68	133.82	122.99
1	E	1	ILE	C-N-CA	7.68	133.82	122.99
1	E	230	VAL	N-CA-C	7.59	118.33	110.36
1	E	87	VAL	CA-C-N	-7.43	110.93	122.37
1	E	87	VAL	C-N-CA	-7.43	110.93	122.37
1	E	268	TYR	CA-C-O	7.35	128.34	120.55
1	E	172	PHE	CA-CB-CG	-7.29	106.51	113.80
1	E	239	LYS	N-CA-CB	7.19	120.69	110.12
1	E	114	PHE	CA-CB-CG	-7.11	106.69	113.80
1	E	29	TYR	N-CA-C	7.10	121.08	109.72
1	E	24	TYR	N-CA-C	7.08	120.06	108.52
1	E	266	ILE	CB-CA-C	-7.04	102.86	111.88
1	E	46	TYR	CB-CA-C	-7.04	100.29	111.48
1	E	283	GLN	N-CA-C	-7.01	103.64	111.28
1	E	67	ASP	CA-CB-CG	7.00	119.60	112.60
1	E	40	PHE	CA-CB-CG	-6.98	106.82	113.80
1	E	114	PHE	N-CA-C	6.90	118.64	108.14
1	E	225	GLN	N-CA-CB	6.89	119.67	110.29
1	E	233	ASN	CA-CB-CG	-6.89	105.71	112.60
1	E	190	GLU	O-C-N	6.83	129.20	122.09
1	E	312	ALA	N-CA-C	6.66	120.67	112.54
1	E	267	PHE	CA-CB-CG	-6.64	107.16	113.80
1	E	261	ASP	N-CA-C	6.61	119.08	111.02
1	E	253	VAL	CB-CA-C	6.60	119.10	110.91
1	E	42	TYR	CB-CA-C	-6.57	97.34	109.37
1	E	122	TYR	N-CA-C	6.55	119.08	108.41
1	E	147	ALA	CA-C-N	-6.53	112.17	120.60
1	E	147	ALA	C-N-CA	-6.53	112.17	120.60
1	E	45	LYS	N-CA-CB	-6.53	102.04	111.70
1	E	24	TYR	CA-CB-CG	-6.51	102.18	113.90
1	E	55	TRP	N-CA-C	6.43	119.95	110.30
1	E	115	TRP	CA-C-N	-6.40	112.43	122.73
1	E	115	TRP	C-N-CA	-6.40	112.43	122.73
1	E	194	THR	CB-CA-C	-6.32	97.71	110.17
1	E	193	TYR	CA-CB-CG	-6.29	102.58	113.90
1	E	268	TYR	N-CA-C	-6.28	104.43	111.28
1	E	19	ASN	N-CA-CB	6.26	120.32	110.56
1	E	24	TYR	N-CA-CB	-6.25	100.73	110.55
1	E	268	TYR	CA-C-N	-6.25	109.90	120.58
1	E	268	TYR	C-N-CA	-6.25	109.90	120.58
1	E	70	ALA	CA-C-O	6.24	127.37	120.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	268	TYR	CA-CB-CG	-6.21	102.72	113.90
1	E	32	ASP	CA-CB-CG	6.16	118.76	112.60
1	E	10	GLY	CA-C-N	-6.14	113.72	122.94
1	E	10	GLY	C-N-CA	-6.14	113.72	122.94
1	E	146	HIS	CA-CB-CG	6.13	119.93	113.80
1	E	193	TYR	N-CA-CB	-6.08	100.38	110.23
1	E	222	THR	O-C-N	6.05	128.83	122.23
1	E	226	ASP	CA-CB-CG	5.97	118.57	112.60
1	E	271	LEU	N-CA-CB	-5.95	101.35	110.16
1	E	239	LYS	CB-CG-CD	5.86	124.77	111.30
1	E	37	ASP	CA-CB-CG	5.82	118.42	112.60
1	E	160	GLU	N-CA-CB	5.81	118.60	109.94
1	E	97	ASN	CA-CB-CG	-5.80	106.80	112.60
1	E	129	THR	CB-CA-C	-5.79	100.10	110.70
1	E	1	ILE	CA-CB-CG2	-5.78	100.68	110.50
1	E	7	VAL	O-C-N	5.78	129.19	123.18
1	E	288	ALA	CA-C-N	-5.76	110.85	120.30
1	E	288	ALA	C-N-CA	-5.76	110.85	120.30
1	E	20	ILE	O-C-N	5.68	127.92	122.97
1	E	73	ALA	CA-C-O	5.67	126.77	120.82
1	E	95	GLY	N-CA-C	-5.67	107.92	115.40
1	E	113	ALA	O-C-N	5.66	129.73	123.33
1	E	16	ASP	CA-CB-CG	5.64	118.24	112.60
1	E	49	THR	O-C-N	-5.62	116.23	122.86
1	E	186	TRP	N-CA-CB	5.58	119.72	110.91
1	E	250	HIS	CA-CB-CG	-5.57	108.23	113.80
1	E	253	VAL	N-CA-CB	5.55	116.78	110.72
1	E	182	LYS	CB-CA-C	5.54	119.34	109.65
1	E	41	THR	CB-CA-C	5.52	119.33	110.16
1	E	279	SER	CB-CA-C	5.52	118.33	109.61
1	E	182	LYS	CG-CD-CE	5.51	123.97	111.30
1	E	256	VAL	N-CA-CB	-5.50	104.73	111.00
1	E	270	ALA	CA-C-O	5.48	126.30	120.10
1	E	242	TYR	N-CA-CB	-5.42	101.93	110.06
1	E	51	PRO	CB-CA-C	-5.42	98.46	112.00
1	E	45	LYS	CB-CA-C	-5.41	104.20	111.89
1	E	269	ARG	O-C-N	-5.38	115.23	122.23
1	E	235	GLY	CA-C-N	5.38	127.34	120.56
1	E	235	GLY	C-N-CA	5.38	127.34	120.56
1	E	6	THR	CA-C-N	-5.38	115.35	122.98
1	E	6	THR	C-N-CA	-5.38	115.35	122.98
1	E	120	MET	CA-C-O	5.37	126.28	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	242	TYR	CA-CB-CG	-5.37	104.24	113.90
1	E	269	ARG	CA-C-O	5.37	126.20	120.24
1	E	316	LYS	CB-CA-C	5.30	120.17	110.10
1	E	20	ILE	N-CA-CB	-5.30	103.26	111.05
1	E	246	GLN	N-CA-C	5.29	119.88	113.38
1	E	170	ASP	CA-C-N	-5.28	114.10	120.70
1	E	170	ASP	C-N-CA	-5.28	114.10	120.70
1	E	188	ILE	O-C-N	5.26	128.71	123.18
1	E	79	VAL	CA-C-O	5.24	126.50	121.05
1	E	59	ASP	N-CA-C	5.24	119.42	112.30
1	E	105	HIS	CA-CB-CG	-5.24	108.56	113.80
1	E	91	LEU	CB-CA-C	-5.23	102.52	111.31
1	E	217	TYR	CA-CB-CG	-5.23	104.49	113.90
1	E	124	ASP	CA-CB-CG	5.21	117.81	112.60
1	E	249	THR	CA-C-O	5.21	125.87	120.40
1	E	281	PHE	N-CA-C	5.20	116.64	111.07
1	E	311	ASP	N-CA-C	-5.20	105.04	111.33
1	E	256	VAL	CA-C-N	5.19	125.93	120.43
1	E	256	VAL	C-N-CA	5.19	125.93	120.43
1	E	178	PHE	CA-CB-CG	-5.17	108.64	113.80
1	E	29	TYR	CB-CA-C	-5.16	99.17	109.33
1	E	20	ILE	N-CA-C	5.16	115.00	107.88
1	E	138	ASP	CA-C-O	5.16	125.93	120.10
1	E	27	TYR	CB-CA-C	-5.15	100.82	111.22
1	E	42	TYR	CA-CB-CG	-5.14	104.65	113.90
1	E	4	THR	N-CA-C	5.11	117.88	109.76
1	E	213	ASP	CB-CA-C	5.10	116.42	108.61
1	E	40	PHE	O-C-N	5.09	129.27	123.31
1	E	274	TYR	N-CA-C	5.09	121.02	114.56
1	E	274	TYR	CB-CA-C	-5.08	101.21	109.13
1	E	76	TYR	CA-CB-CG	-5.07	104.78	113.90
1	E	59	ASP	CA-C-N	-5.06	115.86	124.31
1	E	59	ASP	C-N-CA	-5.06	115.86	124.31
1	E	88	HIS	CA-CB-CG	-5.06	108.74	113.80
1	E	283	GLN	CB-CG-CD	5.05	121.19	112.60
1	E	156	ILE	N-CA-C	5.03	115.56	108.36
1	E	228	GLY	N-CA-C	-5.03	108.30	115.64
1	E	137	ILE	N-CA-C	5.01	116.33	110.62

There are no chirality outliers.

There are no planarity outliers.

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	E	2427	0	2251	222	0
2	E	4	0	0	0	0
3	E	3	0	0	0	0
4	E	127	0	0	10	1
All	All	2561	0	2251	222	1

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 48.

All (222) 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:E:11:ARG:HD2	1:E:63:PHE:CD1	2.03	0.93
1:E:68:ALA:HB3	1:E:69:PRO:HD3	1.48	0.93
1:E:245:SER:HB3	1:E:246:GLN:HG2	1.52	0.89
1:E:11:ARG:HD2	1:E:63:PHE:HD1	1.45	0.82
1:E:262:LYS:O	1:E:266:ILE:HG12	1.81	0.81
1:E:271:LEU:HA	1:E:275:LEU:HD12	1.63	0.79
1:E:1:ILE:HG13	1:E:2:THR:N	2.00	0.77
1:E:57:ASP:HA	4:E:477:HOH:O	1.86	0.76
1:E:1:ILE:HG23	1:E:31:GLN:HE22	1.52	0.74
1:E:30:LEU:HB3	1:E:41:THR:HB	1.70	0.74
1:E:30:LEU:CD1	1:E:70:ALA:HB1	2.18	0.74
1:E:30:LEU:HG	1:E:55:TRP:HB3	1.71	0.73
1:E:240:ALA:O	1:E:244:ILE:HG13	1.89	0.72
1:E:265:LYS:HE3	4:E:403:HOH:O	1.88	0.72
1:E:128:GLN:O	1:E:195:PRO:HD2	1.89	0.71
1:E:42:TYR:CG	1:E:51:PRO:HB2	2.25	0.71
1:E:253:VAL:HG22	1:E:312:ALA:HB2	1.71	0.70
1:E:158:GLN:HG3	1:E:159:ASN:ND2	2.05	0.70
1:E:1:ILE:HG23	1:E:31:GLN:NE2	2.06	0.70
1:E:72:ASP:OD2	1:E:135:GLY:HA2	1.91	0.70
1:E:126:ASP:OD1	1:E:128:GLN:N	2.23	0.70
1:E:167:ALA:O	1:E:171:ILE:HG13	1.92	0.69
1:E:44:ALA:O	1:E:47:ARG:HB2	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:142:HIS:CG	1:E:169:SER:HB3	2.29	0.68
1:E:35:ARG:HD3	1:E:81:TYR:CD1	2.28	0.68
1:E:183:ASN:ND2	4:E:353:HOH:O	2.27	0.68
1:E:164:ILE:O	1:E:168:ILE:HG12	1.94	0.68
1:E:129:THR:HG22	1:E:130:PHE:CD2	2.29	0.67
1:E:30:LEU:HD12	1:E:70:ALA:HB1	1.76	0.67
1:E:301:GLN:O	1:E:304:ALA:HB3	1.94	0.67
1:E:221:TYR:CZ	1:E:223:GLY:HA3	2.30	0.66
1:E:177:GLU:HG2	1:E:184:PRO:HA	1.77	0.65
1:E:37:ASP:HB2	1:E:97:ASN:O	1.96	0.64
1:E:108:GLN:HA	4:E:820:HOH:O	1.97	0.64
1:E:202:LEU:N	1:E:202:LEU:HD22	2.11	0.64
1:E:290:GLN:NE2	1:E:294:ASP:OD1	2.29	0.64
1:E:116:ASN:OD1	1:E:119:GLU:N	2.29	0.64
1:E:182:LYS:C	1:E:184:PRO:HD3	2.21	0.64
1:E:94:ASP:HA	1:E:151:TYR:CD1	2.33	0.64
1:E:68:ALA:CB	1:E:69:PRO:HD3	2.24	0.63
1:E:147:ALA:O	1:E:148:VAL:C	2.41	0.63
1:E:36:GLY:HA3	1:E:97:ASN:ND2	2.14	0.63
1:E:183:ASN:N	1:E:184:PRO:HD3	2.14	0.63
1:E:182:LYS:O	1:E:183:ASN:ND2	2.33	0.61
1:E:217:TYR:O	1:E:220:ARG:HB3	1.99	0.61
1:E:214:PRO:HD2	1:E:232:ILE:CG2	2.31	0.61
1:E:30:LEU:HD23	1:E:30:LEU:N	2.09	0.60
1:E:190:GLU:HG2	1:E:191:ASP:OD1	2.00	0.60
1:E:239:LYS:NZ	4:E:524:HOH:O	2.23	0.60
1:E:4:THR:O	1:E:6:THR:HG23	2.02	0.60
1:E:295:LEU:HB2	1:E:296:TYR:CD2	2.35	0.60
1:E:64:ALA:HB3	1:E:67:ASP:CG	2.26	0.60
1:E:115:TRP:CZ3	1:E:119:GLU:HA	2.36	0.60
1:E:253:VAL:CG2	1:E:312:ALA:HB2	2.31	0.59
1:E:296:TYR:O	1:E:300:SER:HB3	2.02	0.59
1:E:294:ASP:OD1	1:E:294:ASP:N	2.34	0.59
1:E:279:SER:HA	1:E:283:GLN:OE1	2.02	0.59
1:E:188:ILE:O	1:E:201:SER:HB2	2.03	0.59
1:E:40:PHE:CD1	1:E:40:PHE:N	2.69	0.59
1:E:189:GLY:HA3	1:E:202:LEU:HD23	1.85	0.58
1:E:183:ASN:HD22	1:E:183:ASN:C	2.12	0.58
1:E:136:GLY:O	1:E:139:VAL:HB	2.03	0.58
1:E:8:GLY:O	1:E:19:ASN:HB3	2.03	0.57
1:E:85:LYS:HE2	1:E:86:ASN:OD1	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:186:TRP:HB3	1:E:205:MET:HB2	1.86	0.57
1:E:6:THR:HG1	1:E:22:THR:HG1	1.51	0.57
1:E:232:ILE:HD12	1:E:232:ILE:N	2.19	0.57
1:E:110:TYR:CE2	1:E:112:ASN:HB3	2.40	0.57
1:E:42:TYR:HD1	1:E:102:SER:O	1.87	0.57
1:E:10:GLY:HA2	1:E:63:PHE:CE2	2.40	0.56
1:E:188:ILE:O	1:E:202:LEU:HD23	2.05	0.56
1:E:1:ILE:HG23	1:E:1:ILE:O	2.06	0.56
1:E:255:VAL:HG22	1:E:308:GLN:HB2	1.86	0.56
1:E:304:ALA:O	1:E:308:GLN:HG3	2.05	0.56
1:E:271:LEU:HD12	1:E:271:LEU:O	2.06	0.56
1:E:131:ILE:HB	1:E:132:PRO:CD	2.36	0.55
1:E:30:LEU:CG	1:E:55:TRP:HB3	2.35	0.55
1:E:85:LYS:HA	1:E:90:ARG:O	2.06	0.55
1:E:255:VAL:HG13	1:E:308:GLN:OE1	2.06	0.55
1:E:32:ASP:OD2	1:E:35:ARG:NH1	2.38	0.55
1:E:131:ILE:O	1:E:132:PRO:C	2.47	0.55
1:E:19:ASN:C	1:E:20:ILE:HG23	2.33	0.54
1:E:158:GLN:HG3	1:E:159:ASN:HD21	1.72	0.54
1:E:249:THR:HA	1:E:253:VAL:O	2.08	0.54
1:E:190:GLU:HG2	1:E:191:ASP:N	2.22	0.54
1:E:295:LEU:HB2	1:E:296:TYR:CE2	2.43	0.54
1:E:207:ASP:O	1:E:210:LYS:HB3	2.09	0.53
1:E:29:TYR:CD1	1:E:29:TYR:N	2.76	0.53
1:E:70:ALA:O	1:E:74:HIS:HB2	2.08	0.53
1:E:260:ARG:NH1	4:E:513:HOH:O	2.27	0.53
1:E:255:VAL:HA	1:E:308:GLN:OE1	2.08	0.53
1:E:194:THR:OG1	1:E:200:ASP:OD2	2.27	0.53
1:E:19:ASN:O	1:E:20:ILE:HG23	2.09	0.52
1:E:87:VAL:HG12	1:E:88:HIS:CE1	2.44	0.52
1:E:114:PHE:N	1:E:114:PHE:CD1	2.78	0.52
1:E:131:ILE:HB	1:E:132:PRO:HD2	1.91	0.52
1:E:11:ARG:HA	1:E:16:ASP:O	2.09	0.52
1:E:278:THR:O	1:E:279:SER:C	2.53	0.52
1:E:110:TYR:HD2	1:E:123:GLY:HA3	1.75	0.51
1:E:145:THR:O	1:E:148:VAL:HB	2.11	0.51
1:E:68:ALA:C	1:E:71:VAL:HG23	2.36	0.51
1:E:290:GLN:O	1:E:294:ASP:OD1	2.28	0.51
1:E:42:TYR:CD1	1:E:42:TYR:N	2.75	0.50
1:E:50:LEU:HD21	1:E:121:VAL:HG22	1.94	0.50
1:E:244:ILE:O	1:E:258:ILE:HG12	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1:ILE:HG21	1:E:54:LEU:CD2	2.42	0.50
1:E:159:ASN:O	1:E:160:GLU:C	2.54	0.50
1:E:68:ALA:HB3	1:E:69:PRO:CD	2.33	0.50
1:E:151:TYR:N	1:E:151:TYR:CD2	2.79	0.50
1:E:6:THR:HG21	1:E:24:TYR:HB2	1.94	0.49
1:E:29:TYR:HA	1:E:55:TRP:O	2.12	0.49
1:E:310:PHE:O	1:E:311:ASP:C	2.53	0.49
1:E:92:SER:HB2	1:E:97:ASN:HA	1.94	0.49
1:E:305:SER:HA	1:E:308:GLN:CD	2.38	0.49
1:E:240:ALA:HB2	1:E:313:VAL:HG21	1.95	0.48
1:E:261:ASP:O	1:E:265:LYS:HG2	2.14	0.48
1:E:46:TYR:N	1:E:105:HIS:O	2.44	0.48
1:E:68:ALA:CB	1:E:69:PRO:CD	2.90	0.48
1:E:262:LYS:HG3	4:E:528:HOH:O	2.12	0.48
1:E:151:TYR:N	1:E:151:TYR:HD2	2.12	0.48
1:E:156:ILE:HG22	1:E:156:ILE:O	2.14	0.47
1:E:184:PRO:O	1:E:185:ASP:HB3	2.14	0.47
1:E:137:ILE:HG23	1:E:138:ASP:N	2.27	0.47
1:E:253:VAL:HG11	1:E:312:ALA:HA	1.95	0.47
1:E:4:THR:O	1:E:23:THR:HA	2.14	0.47
1:E:30:LEU:HB2	1:E:55:TRP:HB3	1.97	0.47
1:E:1:ILE:CG2	1:E:31:GLN:NE2	2.78	0.47
1:E:189:GLY:HA3	1:E:202:LEU:CD2	2.45	0.47
1:E:214:PRO:HD2	1:E:232:ILE:HG23	1.97	0.47
1:E:40:PHE:O	1:E:101:ARG:HA	2.16	0.46
1:E:160:GLU:HB3	1:E:281:PHE:CD1	2.51	0.46
1:E:253:VAL:HG13	1:E:312:ALA:CB	2.45	0.46
1:E:112:ASN:CG	1:E:113:ALA:N	2.73	0.46
1:E:269:ARG:HD3	1:E:291:SER:HB2	1.97	0.46
1:E:256:VAL:O	1:E:305:SER:HB3	2.15	0.46
1:E:269:ARG:HD3	1:E:291:SER:CB	2.45	0.46
1:E:76:TYR:HD1	1:E:76:TYR:HA	1.49	0.46
1:E:142:HIS:CD2	1:E:142:HIS:C	2.93	0.46
1:E:183:ASN:N	1:E:184:PRO:CD	2.78	0.46
1:E:50:LEU:CD2	1:E:121:VAL:HG22	2.45	0.46
1:E:115:TRP:CZ3	1:E:119:GLU:CA	2.99	0.46
1:E:258:ILE:HG21	1:E:302:GLU:HG3	1.97	0.46
1:E:202:LEU:N	1:E:202:LEU:CD2	2.78	0.46
1:E:297:GLY:C	1:E:299:THR:H	2.23	0.45
1:E:100:ILE:HA	4:E:393:HOH:O	2.15	0.45
1:E:159:ASN:OD1	1:E:228:GLY:HA3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:THR:HA	1:E:195:PRO:HD2	1.36	0.45
1:E:214:PRO:HD2	1:E:232:ILE:HG22	1.96	0.45
1:E:202:LEU:HD22	1:E:202:LEU:H	1.80	0.45
1:E:94:ASP:CG	1:E:96:ASN:H	2.24	0.45
1:E:193:TYR:CE1	1:E:194:THR:HG23	2.52	0.45
1:E:185:ASP:OD2	1:E:190:GLU:OE2	2.35	0.45
1:E:72:ASP:HB3	1:E:134:SER:O	2.17	0.45
1:E:110:TYR:CZ	1:E:112:ASN:HB3	2.52	0.45
1:E:156:ILE:O	1:E:161:SER:HB3	2.17	0.45
1:E:210:LYS:HG2	1:E:211:TYR:CE2	2.52	0.45
1:E:42:TYR:HB3	1:E:51:PRO:O	2.17	0.44
1:E:188:ILE:HD12	1:E:188:ILE:HG23	1.68	0.44
1:E:269:ARG:O	1:E:270:ALA:C	2.57	0.44
1:E:10:GLY:HA3	1:E:62:PHE:HB2	2.00	0.44
1:E:116:ASN:N	1:E:119:GLU:O	2.47	0.44
1:E:27:TYR:CD1	1:E:58:ALA:HA	2.52	0.44
1:E:67:ASP:O	1:E:71:VAL:HG23	2.18	0.44
1:E:30:LEU:HD11	1:E:70:ALA:HB1	1.98	0.44
1:E:291:SER:O	1:E:292:ALA:C	2.58	0.43
1:E:32:ASP:OD1	1:E:34:THR:HG23	2.17	0.43
1:E:87:VAL:HG12	1:E:88:HIS:CD2	2.53	0.43
1:E:106:TYR:O	1:E:107:SER:HB3	2.18	0.43
1:E:167:ALA:O	1:E:168:ILE:C	2.61	0.43
1:E:157:TYR:OH	1:E:166:GLU:OE2	2.29	0.43
1:E:289:VAL:O	1:E:290:GLN:C	2.54	0.43
1:E:182:LYS:HE2	1:E:183:ASN:OD1	2.19	0.43
1:E:202:LEU:HD13	1:E:202:LEU:HA	1.56	0.43
1:E:94:ASP:C	1:E:96:ASN:H	2.27	0.43
1:E:94:ASP:OD1	1:E:94:ASP:N	2.49	0.43
1:E:1:ILE:HG21	1:E:54:LEU:HD22	2.01	0.43
1:E:177:GLU:OE2	1:E:184:PRO:HA	2.19	0.43
1:E:57:ASP:OD2	1:E:60:ASN:N	2.52	0.43
1:E:109:GLY:HA2	1:E:125:GLY:O	2.18	0.43
1:E:131:ILE:CB	1:E:132:PRO:CD	2.95	0.43
1:E:193:TYR:CZ	1:E:194:THR:HG23	2.54	0.43
1:E:239:LYS:HE3	1:E:243:LEU:HD11	2.01	0.43
1:E:12:GLY:N	1:E:16:ASP:O	2.45	0.43
1:E:43:ASP:O	1:E:52:GLY:HA3	2.19	0.43
1:E:104:VAL:O	1:E:105:HIS:HB2	2.19	0.43
1:E:146:HIS:ND1	1:E:165:ASN:OD1	2.52	0.43
1:E:175:LEU:HD23	1:E:175:LEU:HA	1.74	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:269:ARG:HB2	4:E:335:HOH:O	2.18	0.43
1:E:295:LEU:CB	1:E:296:TYR:CE2	3.02	0.43
1:E:53:SER:O	1:E:54:LEU:C	2.61	0.42
1:E:85:LYS:O	1:E:89:ASN:HA	2.18	0.42
1:E:110:TYR:HD2	1:E:123:GLY:CA	2.32	0.42
1:E:130:PHE:C	1:E:131:ILE:HG23	2.44	0.42
1:E:157:TYR:CE1	1:E:166:GLU:HG2	2.55	0.42
1:E:289:VAL:CG2	1:E:290:GLN:N	2.79	0.42
1:E:289:VAL:HG12	1:E:306:VAL:HG12	2.01	0.42
1:E:62:PHE:N	1:E:62:PHE:CD1	2.86	0.42
1:E:25:SER:O	1:E:27:TYR:N	2.48	0.42
1:E:243:LEU:HD12	1:E:243:LEU:HA	1.70	0.42
1:E:11:ARG:HE	1:E:17:GLN:HG3	1.83	0.42
1:E:189:GLY:O	1:E:190:GLU:C	2.61	0.42
1:E:217:TYR:HB2	1:E:313:VAL:O	2.20	0.42
1:E:276:THR:C	1:E:278:THR:N	2.77	0.42
1:E:271:LEU:HD12	1:E:271:LEU:C	2.42	0.41
1:E:31:GLN:HG3	1:E:40:PHE:CE2	2.55	0.41
1:E:293:THR:O	1:E:294:ASP:C	2.61	0.41
1:E:315:VAL:HA	4:E:371:HOH:O	2.20	0.41
1:E:94:ASP:C	1:E:96:ASN:N	2.79	0.41
1:E:11:ARG:NE	1:E:17:GLN:HG3	2.36	0.41
1:E:189:GLY:O	1:E:193:TYR:HB2	2.20	0.41
1:E:25:SER:O	1:E:26:THR:HB	2.21	0.41
1:E:194:THR:N	1:E:195:PRO:CD	2.75	0.41
1:E:30:LEU:HD22	1:E:30:LEU:HA	1.75	0.41
1:E:146:HIS:CE1	1:E:166:GLU:OE2	2.72	0.40
1:E:11:ARG:HD2	1:E:63:PHE:CE1	2.55	0.40
1:E:11:ARG:HE	1:E:11:ARG:HB3	1.42	0.40
1:E:115:TRP:CE3	1:E:119:GLU:C	2.99	0.40
1:E:129:THR:CG2	1:E:130:PHE:CD2	3.03	0.40
1:E:44:ALA:O	1:E:45:LYS:HB2	2.21	0.40
1:E:138:ASP:OD2	1:E:188:ILE:HA	2.22	0.40
1:E:67:ASP:C	1:E:69:PRO:HD2	2.47	0.40
1:E:267:PHE:O	1:E:268:TYR:C	2.61	0.40

All (1) 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)
4:E:509:HOH:O	4:E:729:HOH:O[5_545]	2.19	0.01

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	E	314/316 (99%)	286 (91%)	24 (8%)	4 (1%)	9	10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	68	ALA
1	E	218	SER
1	E	132	PRO
1	E	183	ASN

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	E	250/252 (99%)	208 (83%)	42 (17%)	2	2

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	1	ILE
1	E	2	THR
1	E	11	ARG
1	E	16	ASP
1	E	19	ASN
1	E	20	ILE
1	E	25	SER

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Mol	Chain	Res	Type
1	E	30	LEU
1	E	34	THR
1	E	48	THR
1	E	49	THR
1	E	50	LEU
1	E	53	SER
1	E	71	VAL
1	E	85	LYS
1	E	97	ASN
1	E	101	ARG
1	E	102	SER
1	E	108	GLN
1	E	118	SER
1	E	156	ILE
1	E	157	TYR
1	E	159	ASN
1	E	161	SER
1	E	174	THR
1	E	182	LYS
1	E	183	ASN
1	E	198	SER
1	E	220	ARG
1	E	225	GLN
1	E	234	SER
1	E	243	LEU
1	E	253	VAL
1	E	261	ASP
1	E	265	LYS
1	E	273	GLN
1	E	289	VAL
1	E	291	SER
1	E	294	ASP
1	E	303	VAL
1	E	305	SER
1	E	308	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	31	GLN
1	E	33	ASN
1	E	97	ASN

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Mol	Chain	Res	Type
1	E	128	GLN
1	E	158	GLN
1	E	159	ASN
1	E	183	ASN
1	E	227	ASN
1	E	273	GLN

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](#)

Of 7 ligands modelled in this entry, 7 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 [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	E	316/316 (100%)	0.20	4 (1%) 75 76	53, 70, 89, 100	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	157	TYR	2.8
1	E	182	LYS	2.6
1	E	12	GLY	2.5
1	E	16	ASP	2.1

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	CA	E	319	1/1	0.99	0.04	81,81,81,81	0
2	CA	E	320	1/1	0.99	0.04	73,73,73,73	0
3	ZN	E	322	1/1	0.99	0.06	64,64,64,64	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	E	330	1/1	0.99	0.03	70,70,70,70	0
3	ZN	E	321	1/1	1.00	0.02	62,62,62,62	0
2	CA	E	317	1/1	1.00	0.02	59,59,59,59	0
2	CA	E	318	1/1	1.00	0.02	53,53,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.