



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:56 PM UTC

PDB ID : 4IBX / pdb_00004ibx
Title : Crystal structure of stabilized TEM-1 beta-lactamase variant v.13
Authors : Dellus-Gur, E.; Toth-Petroczy, A.; Elias, M.; Tawfik, D.S.
Deposited on : 2012-12-09
Resolution : 2.68 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

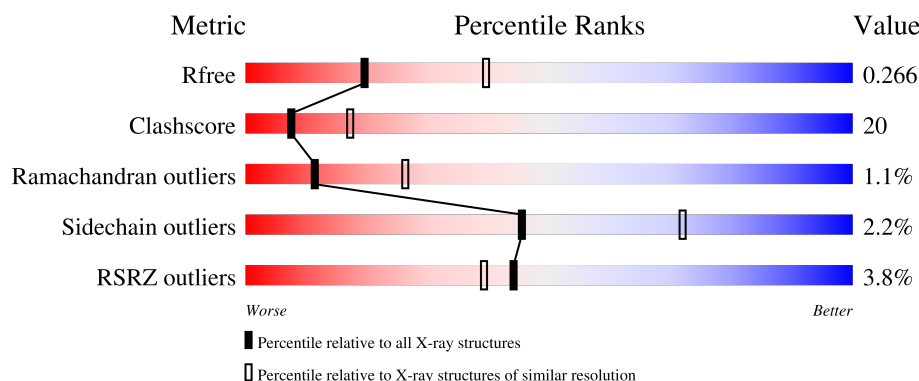
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	263	71% 26% .
1	B	263	70% 28% .
1	C	263	2% 59% 38% .
1	D	263	3% 67% 32%
1	E	263	12% 56% 40% 5%

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	CA	A	301	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10281 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 Beta-lactamase TEM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	2	0
			2028	1267	358	392	11			
1	B	263	Total	C	N	O	S	0	2	0
			2030	1270	356	393	11			
1	C	263	Total	C	N	O	S	0	3	0
			2033	1271	358	393	11			
1	D	263	Total	C	N	O	S	0	2	0
			2028	1267	359	391	11			
1	E	263	Total	C	N	O	S	0	1	0
			2023	1263	359	390	11			

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	GLY	ALA	engineered mutation	UNP P62593
A	52	ALA	ASN	engineered mutation	UNP P62593
A	120	GLY	ARG	engineered mutation	UNP P62593
A	182	THR	MET	engineered mutation	UNP P62593
A	184	VAL	ALA	conflict	UNP P62593
A	201	ALA	LEU	engineered mutation	UNP P62593
A	265	MET	THR	engineered mutation	UNP P62593
B	42	GLY	ALA	engineered mutation	UNP P62593
B	52	ALA	ASN	engineered mutation	UNP P62593
B	120	GLY	ARG	engineered mutation	UNP P62593
B	182	THR	MET	engineered mutation	UNP P62593
B	184	VAL	ALA	conflict	UNP P62593
B	201	ALA	LEU	engineered mutation	UNP P62593
B	265	MET	THR	engineered mutation	UNP P62593
C	42	GLY	ALA	engineered mutation	UNP P62593
C	52	ALA	ASN	engineered mutation	UNP P62593
C	120	GLY	ARG	engineered mutation	UNP P62593
C	182	THR	MET	engineered mutation	UNP P62593
C	184	VAL	ALA	conflict	UNP P62593

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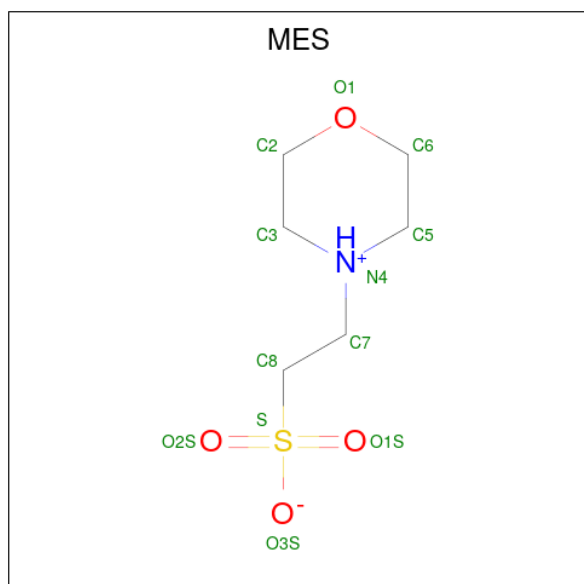
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Chain	Residue	Modelled	Actual	Comment	Reference
C	201	ALA	LEU	engineered mutation	UNP P62593
C	265	MET	THR	engineered mutation	UNP P62593
D	42	GLY	ALA	engineered mutation	UNP P62593
D	52	ALA	ASN	engineered mutation	UNP P62593
D	120	GLY	ARG	engineered mutation	UNP P62593
D	182	THR	MET	engineered mutation	UNP P62593
D	184	VAL	ALA	conflict	UNP P62593
D	201	ALA	LEU	engineered mutation	UNP P62593
D	265	MET	THR	engineered mutation	UNP P62593
E	42	GLY	ALA	engineered mutation	UNP P62593
E	52	ALA	ASN	engineered mutation	UNP P62593
E	120	GLY	ARG	engineered mutation	UNP P62593
E	182	THR	MET	engineered mutation	UNP P62593
E	184	VAL	ALA	conflict	UNP P62593
E	201	ALA	LEU	engineered mutation	UNP P62593
E	265	MET	THR	engineered mutation	UNP P62593

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

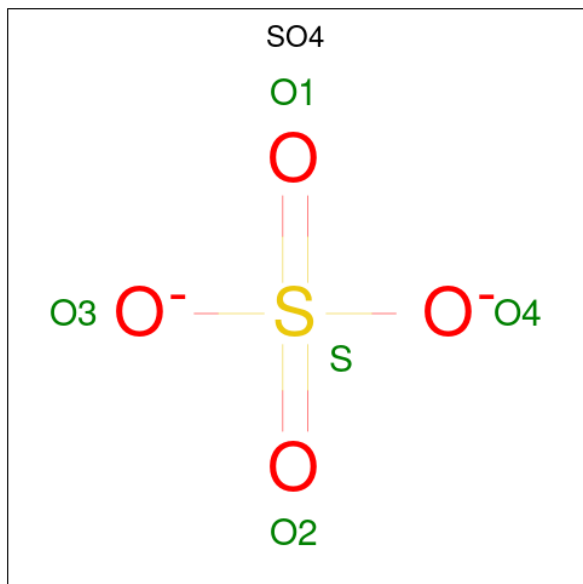
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0

- Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
3	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

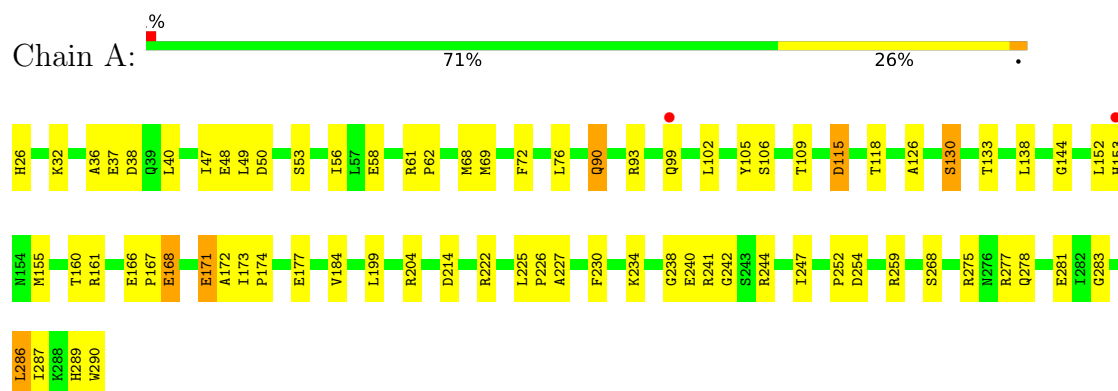
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	34	Total	O	0	0
			34	34		
5	B	30	Total	O	0	0
			30	30		
5	C	11	Total	O	0	0
			11	11		
5	D	10	Total	O	0	0
			10	10		
5	E	11	Total	O	0	0
			11	11		

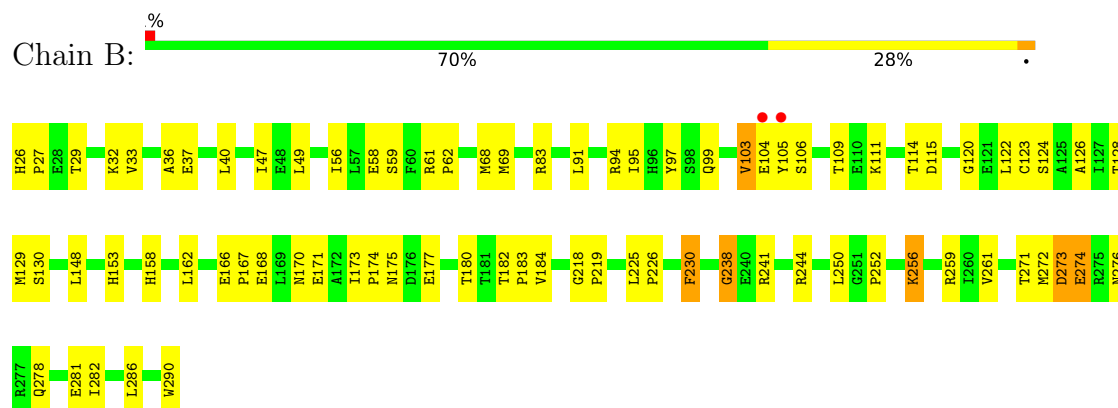
3 Residue-property plots

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.

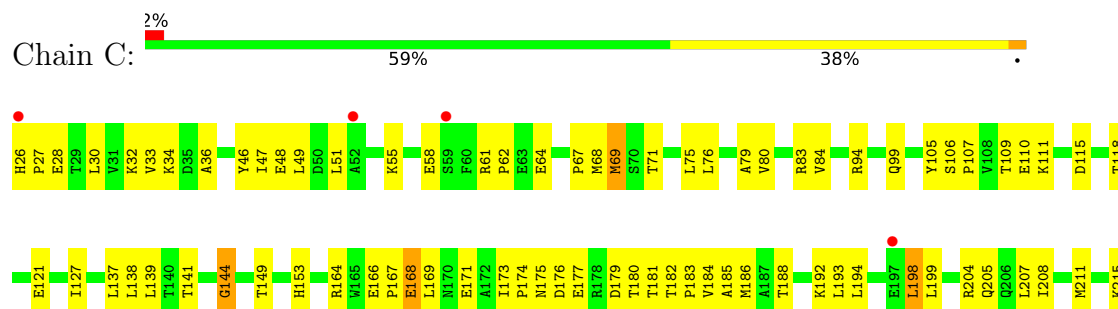
• Molecule 1: Beta-lactamase TEM



• Molecule 1: Beta-lactamase TEM

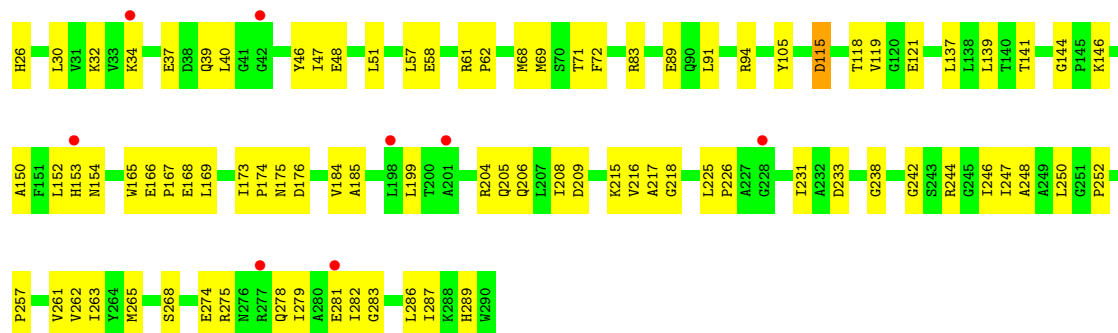


• Molecule 1: Beta-lactamase TEM

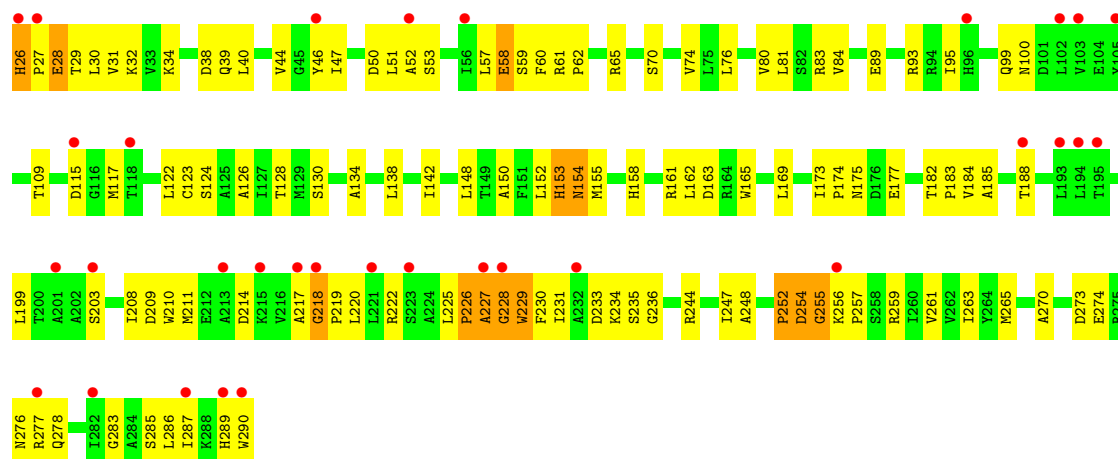




● Molecule 1: Beta-lactamase TEM



● Molecule 1: Beta-lactamase TEM



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	39.90Å 239.00Å 159.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.68 – 2.68 48.68 – 2.68	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.68-2.68) 98.4 (48.68-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.232 , 0.265 0.232 , 0.266	Depositor DCC
R_{free} test set	2187 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10281	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.89 % 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 2.2090e-04. 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, SO4, MES

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	0.45	1/2069 (0.0%)	0.81	7/2801 (0.2%)
1	B	0.51	2/2071 (0.1%)	0.77	7/2804 (0.2%)
1	C	0.46	2/2076 (0.1%)	0.73	3/2809 (0.1%)
1	D	0.51	1/2069 (0.0%)	0.81	4/2801 (0.1%)
1	E	0.41	1/2060 (0.0%)	0.81	5/2788 (0.2%)
All	All	0.47	7/10345 (0.1%)	0.79	26/14003 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	252	PRO	C-N	16.60	1.56	1.33
1	B	252	PRO	C-N	16.12	1.53	1.33
1	C	252	PRO	C-N	13.42	1.52	1.33
1	A	252	PRO	C-N	12.19	1.50	1.33
1	E	252	PRO	C-N	10.37	1.48	1.33
1	C	238	GLY	C-N	8.11	1.44	1.33
1	B	256	LYS	C-O	-5.49	1.16	1.24

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	252	PRO	CA-C-N	-9.80	102.82	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	252	PRO	C-N-CA	-9.80	102.82	121.54
1	A	238	GLY	O-C-N	9.77	132.33	123.73
1	E	153	HIS	N-CA-C	-8.80	97.57	110.42
1	D	238	GLY	O-C-N	8.30	133.49	122.70
1	D	252	PRO	CA-C-N	-7.64	111.74	122.74
1	D	252	PRO	C-N-CA	-7.64	111.74	122.74
1	B	238	GLY	CA-C-N	-7.17	108.46	122.02
1	B	238	GLY	C-N-CA	-7.17	108.46	122.02
1	A	238	GLY	CA-C-N	-7.14	107.67	122.03
1	A	238	GLY	C-N-CA	-7.14	107.67	122.03
1	B	274	GLU	N-CA-C	-7.08	103.31	112.23
1	B	252	PRO	CA-C-N	-7.03	112.66	125.02
1	B	252	PRO	C-N-CA	-7.03	112.66	125.02
1	D	252	PRO	O-C-N	6.69	130.66	123.23
1	C	144	GLY	CA-C-N	6.59	125.82	118.97
1	C	144	GLY	C-N-CA	6.59	125.82	118.97
1	A	144	GLY	CA-C-N	6.04	127.39	119.84
1	A	144	GLY	C-N-CA	6.04	127.39	119.84
1	E	26	HIS	CA-C-N	5.90	127.21	119.84
1	E	26	HIS	C-N-CA	5.90	127.21	119.84
1	C	252	PRO	O-C-N	-5.72	114.92	122.64
1	A	252	PRO	CA-C-N	-5.54	114.99	123.47
1	A	252	PRO	C-N-CA	-5.54	114.99	123.47
1	B	252	PRO	O-C-N	5.07	129.12	123.04
1	B	238	GLY	O-C-N	5.06	129.28	122.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	252	PRO	Mainchain

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	2028	0	2032	57	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2030	0	2034	63	0
1	C	2033	0	2046	95	1
1	D	2028	0	2034	67	0
1	E	2023	0	2032	135	2
2	A	1	0	0	2	0
2	C	1	0	0	0	0
3	A	12	0	12	0	0
3	B	12	0	12	2	0
3	D	12	0	12	2	0
4	C	5	0	0	1	0
5	A	34	0	0	4	0
5	B	30	0	0	1	0
5	C	11	0	0	4	0
5	D	10	0	0	3	0
5	E	11	0	0	0	0
All	All	10281	0	10214	415	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 20.

All (415) 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:230:PHE:CE1	1:E:257:PRO:HD3	1.23	1.74
1:E:230:PHE:CD1	1:E:257:PRO:HD3	1.37	1.54
1:E:26:HIS:CE1	1:E:289:HIS:ND1	1.72	1.54
1:E:244:ARG:HB3	1:E:265:MET:CE	1.54	1.34
1:E:230:PHE:CD1	1:E:257:PRO:CD	2.13	1.30
1:E:230:PHE:CE1	1:E:257:PRO:CD	2.16	1.29
1:E:26:HIS:ND1	1:E:289:HIS:CG	2.02	1.26
1:C:68:MET:HE3	1:C:186:MET:HE1	1.27	1.11
1:E:26:HIS:ND1	1:E:289:HIS:HB3	1.65	1.11
1:B:273:ASP:HA	1:B:276:ASN:HB2	1.22	1.10
1:C:68:MET:HE3	1:C:186:MET:CE	1.83	1.08
1:E:252:PRO:O	1:E:254:ASP:HB2	1.50	1.08
1:E:26:HIS:ND1	1:E:289:HIS:CB	2.18	1.07
1:E:29:THR:OG1	1:E:285:SER:O	1.70	1.06
1:E:26:HIS:HE1	1:E:289:HIS:ND1	1.21	1.04
1:E:244:ARG:CB	1:E:265:MET:CE	2.36	1.04
1:B:271:THR:OG1	1:B:274:GLU:HG3	1.54	1.03
1:E:153:HIS:O	1:E:154:ASN:CB	2.09	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:153:HIS:O	1:E:154:ASN:HB2	1.19	1.00
1:E:244:ARG:CB	1:E:265:MET:HE2	1.91	1.00
1:E:26:HIS:CE1	1:E:289:HIS:CE1	2.51	0.99
1:E:230:PHE:HD1	1:E:257:PRO:N	1.62	0.98
1:E:26:HIS:ND1	1:E:289:HIS:ND1	2.05	0.97
1:E:26:HIS:HE1	1:E:289:HIS:CE1	1.81	0.97
1:C:205:GLN:NE2	1:C:208:ILE:HD11	1.80	0.95
1:E:244:ARG:HB3	1:E:265:MET:HE2	0.96	0.95
1:E:30:LEU:HD22	1:E:58:GLU:OE1	1.69	0.93
1:B:103:VAL:HB	1:B:104[B]:GLU:OE1	1.68	0.93
1:E:26:HIS:HD2	1:E:27:PRO:HD2	1.34	0.91
1:C:252:PRO:O	1:C:254:ASP:HB2	1.69	0.90
1:C:205:GLN:HE21	1:C:208:ILE:HD11	1.34	0.89
1:C:68:MET:CE	1:C:186:MET:HE1	2.03	0.89
1:E:26:HIS:CE1	1:E:289:HIS:CG	2.53	0.88
1:E:228:GLY:O	1:E:229:TRP:HB2	1.74	0.87
1:C:34[A]:LYS:HG2	5:C:407:HOH:O	1.73	0.87
1:E:230:PHE:HE1	1:E:257:PRO:HD3	1.08	0.87
1:E:26:HIS:CD2	1:E:27:PRO:HD2	2.11	0.86
1:E:230:PHE:HD1	1:E:257:PRO:CD	1.70	0.86
1:C:32:LYS:HE2	1:C:281:GLU:HB3	1.58	0.84
1:E:30:LEU:HD12	1:E:30:LEU:O	1.79	0.82
1:E:225:LEU:HD22	1:E:231:ILE:HD12	1.61	0.82
1:E:99:GLN:NE2	1:E:100:ASN:OD1	2.12	0.82
1:E:30:LEU:CD2	1:E:58:GLU:OE1	2.27	0.81
1:E:214:ASP:HB3	1:E:222:ARG:HH21	1.46	0.81
1:C:215:LYS:HE2	1:C:215:LYS:HA	1.63	0.81
1:E:254:ASP:O	1:E:256:LYS:N	2.13	0.81
1:E:230:PHE:HE1	1:E:257:PRO:CD	1.75	0.80
1:C:198:LEU:HD22	1:C:198:LEU:H	1.45	0.79
1:B:61:ARG:N	1:B:62:PRO:HD3	1.96	0.79
1:B:273:ASP:HA	1:B:276:ASN:CB	2.09	0.79
1:D:40:LEU:HD11	1:D:278:GLN:HG3	1.63	0.78
1:D:62:PRO:O	1:D:184:VAL:HG12	1.83	0.77
1:A:38:ASP:OD1	5:A:425:HOH:O	2.01	0.77
1:C:194:LEU:HD22	1:C:208:ILE:HG23	1.65	0.77
1:B:244:ARG:NH1	3:B:301:MES:O3S	2.19	0.75
1:E:230:PHE:CD1	1:E:257:PRO:N	2.44	0.75
1:C:68:MET:HE1	1:C:75:LEU:HD12	1.70	0.74
1:B:103:VAL:C	1:B:104[B]:GLU:OE1	2.31	0.74
1:D:89:GLU:OE2	5:D:407:HOH:O	2.05	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:GLY:O	1:C:287:ILE:HG12	1.87	0.73
1:D:37:GLU:OE2	1:D:61:ARG:NH2	2.20	0.73
1:D:268:SER:HB2	5:D:406:HOH:O	1.88	0.73
1:E:32:LYS:O	1:E:32:LYS:HD3	1.87	0.73
1:B:103:VAL:CB	1:B:104[B]:GLU:OE1	2.38	0.72
1:C:244:ARG:NH1	4:C:302:SO4:O2	2.22	0.71
1:C:32:LYS:HE2	1:C:281:GLU:CB	2.20	0.71
1:B:59:SER:HB2	1:B:62:PRO:HG3	1.72	0.71
1:E:230:PHE:HD1	1:E:256:LYS:C	1.97	0.71
1:C:79:ALA:O	1:C:83:ARG:HG3	1.91	0.71
1:E:230:PHE:CD1	1:E:256:LYS:HA	2.26	0.70
1:E:150:ALA:O	1:E:153:HIS:O	2.08	0.70
1:E:231:ILE:HG23	1:E:231:ILE:O	1.89	0.70
1:D:283:GLY:O	1:D:287:ILE:HG12	1.92	0.70
1:E:26:HIS:CG	1:E:289:HIS:HB3	2.26	0.70
1:C:46:TYR:HD1	1:C:263:ILE:HG12	1.56	0.69
1:E:30:LEU:HB2	1:E:58:GLU:OE1	1.92	0.69
1:B:230:PHE:C	1:B:230:PHE:CD1	2.71	0.69
1:D:215:LYS:HE2	1:D:215:LYS:HA	1.75	0.69
1:E:46:TYR:HD1	1:E:263:ILE:HG12	1.59	0.68
1:E:153:HIS:ND1	1:E:158:HIS:HD2	1.92	0.67
1:B:272:MET:O	1:B:273:ASP:CB	2.42	0.67
1:E:29:THR:OG1	1:E:285:SER:OG	2.06	0.67
1:D:40:LEU:CD1	1:D:278:GLN:HG3	2.25	0.66
1:A:171:GLU:OE2	1:A:173:ILE:HD11	1.96	0.65
1:E:29:THR:OG1	1:E:285:SER:C	2.40	0.65
1:A:268:SER:HB3	5:A:418:HOH:O	1.96	0.65
1:C:106:SER:HB3	1:C:109:THR:OG1	1.97	0.65
1:D:32:LYS:HE2	1:D:281:GLU:HB2	1.80	0.64
1:E:225:LEU:O	1:E:226:PRO:O	2.15	0.64
1:E:244:ARG:HB3	1:E:265:MET:HE3	1.72	0.64
1:C:47:ILE:HG22	1:C:262:VAL:HB	1.80	0.63
1:A:286:LEU:CD1	1:A:290:TRP:HZ3	2.12	0.63
1:E:59:SER:HB2	1:E:62:PRO:HG3	1.79	0.63
1:E:254:ASP:O	1:E:255:GLY:C	2.42	0.63
1:A:26:HIS:CD2	1:A:289:HIS:HB3	2.33	0.63
1:B:26:HIS:N	1:B:58:GLU:HG3	2.12	0.63
1:C:76:LEU:HD21	1:C:138:LEU:HB2	1.79	0.63
1:E:109:THR:HG22	1:E:117:MET:SD	2.38	0.63
1:B:272:MET:O	1:B:273:ASP:HB3	1.97	0.62
1:B:103:VAL:CA	1:B:104[B]:GLU:OE1	2.47	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:MET:O	1:B:273:ASP:OD2	2.16	0.62
1:E:230:PHE:CE1	1:E:257:PRO:CG	2.81	0.62
2:A:301:CA:CA	1:B:37:GLU:OE1	1.77	0.62
1:E:228:GLY:O	1:E:229:TRP:CB	2.47	0.62
1:D:47:ILE:HG22	1:D:262:VAL:HB	1.81	0.62
1:D:247:ILE:HG13	1:D:247:ILE:O	1.99	0.62
1:D:265:MET:CE	1:D:279:ILE:HD11	2.30	0.61
1:C:198:LEU:H	1:C:198:LEU:CD2	2.13	0.61
1:E:84:VAL:HA	1:E:89:GLU:O	2.00	0.61
1:C:68:MET:HG2	1:C:181:THR:HG22	1.83	0.61
1:A:37:GLU:OE1	2:A:301:CA:CA	1.79	0.60
1:D:61:ARG:N	1:D:62:PRO:CD	2.63	0.60
1:E:222:ARG:HG2	1:E:231:ILE:HD13	1.83	0.60
1:E:199:LEU:HB3	1:E:203:SER:HB2	1.83	0.60
1:A:230:PHE:CD1	1:A:230:PHE:C	2.79	0.59
1:C:164:ARG:NE	1:C:179:ASP:OD2	2.32	0.59
1:C:198:LEU:HD22	1:C:198:LEU:N	2.16	0.59
1:E:169:LEU:C	1:E:169:LEU:HD12	2.27	0.59
1:E:26:HIS:CE1	1:E:28:GLU:HG3	2.37	0.59
1:E:261:VAL:HG21	1:E:286:LEU:HD21	1.83	0.59
1:C:167:PRO:O	1:C:169:LEU:N	2.36	0.59
1:A:161:ARG:HD3	1:A:177:GLU:O	2.03	0.58
1:E:57:LEU:O	1:E:58:GLU:HB2	2.02	0.58
1:C:80:VAL:O	1:C:84:VAL:HG23	2.03	0.58
1:E:231:ILE:O	1:E:231:ILE:CG2	2.51	0.58
1:B:170:ASN:HB3	1:B:238:GLY:HA2	1.84	0.58
1:A:214:ASP:HB3	1:A:222:ARG:HH21	1.68	0.58
1:B:95:ILE:HD12	1:B:122:LEU:HD12	1.85	0.57
1:A:99:GLN:OE1	1:A:99:GLN:N	2.24	0.57
1:E:230:PHE:HD1	1:E:256:LYS:CA	2.17	0.57
1:C:26:HIS:ND1	1:C:27:PRO:HD2	2.19	0.57
1:A:115:ASP:OD2	5:A:420:HOH:O	2.18	0.56
1:E:220:LEU:HD12	1:E:235:SER:HB3	1.86	0.56
1:E:230:PHE:CD1	1:E:256:LYS:CA	2.88	0.56
1:A:99:GLN:H	1:A:99:GLN:CD	2.11	0.56
1:A:286:LEU:HD11	1:A:290:TRP:HZ3	1.70	0.56
1:C:61:ARG:O	1:C:183:PRO:HG2	2.04	0.56
1:A:106:SER:HB3	1:A:109:THR:OG1	2.06	0.56
1:C:250:LEU:HD23	1:C:250:LEU:H	1.71	0.56
1:B:26:HIS:N	1:B:58:GLU:CG	2.69	0.56
1:C:68:MET:O	1:C:71:THR:OG1	2.15	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:TYR:C	1:C:107:PRO:HD3	2.31	0.56
1:D:40:LEU:HD11	1:D:278:GLN:CG	2.35	0.56
1:D:115:ASP:OD1	1:D:115:ASP:N	2.38	0.56
1:C:62:PRO:HA	1:C:183:PRO:HB2	1.88	0.56
1:C:182:THR:OG1	1:C:185:ALA:CB	2.53	0.56
1:B:126:ALA:O	1:B:130:SER:HA	2.06	0.56
1:D:139:LEU:O	1:D:144:GLY:N	2.36	0.56
1:C:36:ALA:HA	1:C:278:GLN:OE1	2.06	0.55
1:C:261:VAL:HG21	1:C:286:LEU:CD2	2.37	0.55
1:E:214:ASP:OD2	1:E:217:ALA:HB2	2.06	0.55
1:C:139:LEU:O	1:C:144:GLY:N	2.35	0.55
1:A:48:GLU:OE2	1:A:259:ARG:NH1	2.37	0.54
1:A:102:LEU:HD23	1:A:133:THR:HG21	1.89	0.54
1:C:68:MET:HG2	1:C:181:THR:CG2	2.37	0.54
1:C:177:GLU:O	1:C:180:THR:OG1	2.15	0.54
1:E:152:LEU:O	1:E:153:HIS:C	2.47	0.54
1:D:250:LEU:H	1:D:250:LEU:HD23	1.72	0.54
1:A:277:ARG:O	1:A:281:GLU:HG3	2.08	0.53
1:A:283:GLY:HA2	1:A:286:LEU:HD23	1.91	0.53
1:A:40:LEU:HD21	1:A:278:GLN:HG3	1.91	0.53
1:E:236:GLY:O	1:E:244:ARG:NH1	2.35	0.53
1:E:76:LEU:HD21	1:E:138:LEU:HB2	1.91	0.53
1:B:83:ARG:NH1	5:B:419:HOH:O	2.42	0.53
1:E:44:VAL:O	1:E:60:PHE:HD2	1.90	0.53
1:D:169:LEU:HD12	1:D:169:LEU:C	2.34	0.53
1:D:47:ILE:CG2	1:D:262:VAL:HB	2.38	0.53
1:B:271:THR:OG1	1:B:274:GLU:CG	2.45	0.53
1:E:225:LEU:HD21	1:E:231:ILE:HB	1.91	0.53
1:B:153:HIS:HD2	1:B:158:HIS:CD2	2.28	0.52
1:C:199:LEU:HB2	1:C:204:ARG:HG3	1.92	0.52
1:C:184:VAL:HG13	1:C:185:ALA:N	2.25	0.52
1:E:247:ILE:HG13	1:E:247:ILE:O	2.08	0.52
1:A:268:SER:CB	5:A:418:HOH:O	2.56	0.52
1:D:174:PRO:O	1:D:175:ASN:HB2	2.08	0.52
1:E:29:THR:HA	1:E:285:SER:OG	2.10	0.52
1:A:50:ASP:HB3	1:A:53:SER:OG	2.09	0.52
1:E:274:GLU:O	1:E:278:GLN:HG2	2.09	0.52
1:D:118:THR:OG1	1:D:121:GLU:HG3	2.10	0.52
1:E:214:ASP:HB3	1:E:222:ARG:NH2	2.21	0.52
1:E:30:LEU:HD11	1:E:34:LYS:HE3	1.92	0.52
1:E:225:LEU:CD2	1:E:231:ILE:HB	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:ARG:N	1:B:62:PRO:CD	2.70	0.51
1:C:205:GLN:HE21	1:C:208:ILE:CD1	2.16	0.51
1:E:44:VAL:O	1:E:60:PHE:CD2	2.64	0.51
1:C:69:MET:HE2	1:C:243:SER:HB3	1.93	0.51
1:E:227:ALA:O	1:E:228:GLY:O	2.28	0.51
1:A:61:ARG:N	1:A:62:PRO:CD	2.73	0.51
1:C:247:ILE:HG13	1:C:247:ILE:O	2.11	0.51
1:B:36:ALA:O	1:B:40:LEU:HG	2.11	0.51
1:B:106:SER:HB3	1:B:109:THR:OG1	2.11	0.51
1:B:261:VAL:HG21	1:B:286:LEU:CD2	2.41	0.51
1:D:274:GLU:O	1:D:278:GLN:HG2	2.11	0.51
1:A:254:ASP:OD1	1:D:141:THR:HG22	2.11	0.50
1:E:259:ARG:HB2	1:E:290:TRP:HH2	1.75	0.50
1:C:251:GLY:HA3	1:C:257:PRO:HA	1.93	0.50
1:B:278:GLN:O	1:B:282:ILE:HG13	2.11	0.50
1:C:64:GLU:O	1:C:183:PRO:HD3	2.12	0.50
1:D:61:ARG:N	1:D:62:PRO:HD3	2.26	0.50
1:E:217:ALA:O	1:E:218:GLY:C	2.54	0.50
1:C:218:GLY:N	1:C:219:PRO:HD2	2.27	0.50
1:C:261:VAL:HG21	1:C:286:LEU:HD21	1.93	0.50
1:D:199:LEU:HB2	1:D:204:ARG:HG3	1.94	0.50
1:E:126:ALA:O	1:E:130:SER:HA	2.12	0.50
1:E:47:ILE:CD1	1:E:184:VAL:HA	2.42	0.49
1:E:153:HIS:ND1	1:E:158:HIS:CD2	2.78	0.49
1:D:34:LYS:O	1:D:34:LYS:HD3	2.11	0.49
1:C:167:PRO:C	1:C:169:LEU:N	2.69	0.49
1:C:188:THR:O	1:C:192:LYS:HG3	2.13	0.49
1:C:182:THR:OG1	1:C:185:ALA:HB2	2.13	0.49
1:C:105:TYR:O	1:C:107:PRO:HD3	2.13	0.49
1:A:40:LEU:HD21	1:A:278:GLN:CG	2.43	0.49
1:A:172:ALA:O	1:A:241:ARG:HD2	2.13	0.49
1:A:286:LEU:HD12	1:A:290:TRP:CZ3	2.48	0.49
1:D:250:LEU:HD23	1:D:250:LEU:N	2.27	0.49
1:E:39:GLN:HE22	1:E:278:GLN:NE2	2.11	0.48
1:E:50:ASP:OD2	1:E:53:SER:OG	2.31	0.48
1:D:46:TYR:HD1	1:D:263:ILE:HG12	1.79	0.48
1:B:95:ILE:HD12	1:B:122:LEU:CD1	2.44	0.48
1:D:278:GLN:O	1:D:282:ILE:HG13	2.12	0.48
1:D:173:ILE:HB	1:D:176:ASP:HB2	1.95	0.48
1:B:271:THR:HG1	1:B:274:GLU:HG3	1.71	0.48
1:E:173:ILE:HG23	1:E:174:PRO:HD2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:261:VAL:HG21	1:E:286:LEU:CD2	2.43	0.48
1:A:68:MET:O	1:A:69:MET:HB2	2.14	0.48
1:B:174:PRO:O	1:B:175:ASN:HB2	2.13	0.48
1:D:216:VAL:HB	3:D:301:MES:C6	2.44	0.48
1:C:230:PHE:CD1	1:C:230:PHE:C	2.91	0.48
1:C:236:GLY:O	1:C:244:ARG:HD2	2.14	0.48
1:E:40:LEU:HD22	1:E:270:ALA:HB3	1.95	0.48
1:B:124:SER:O	1:B:128:THR:OG1	2.21	0.48
1:D:165:TRP:CD1	1:D:165:TRP:N	2.82	0.48
1:A:166:GLU:OE1	1:A:167:PRO:HA	2.15	0.47
1:B:105[A]:TYR:HE1	1:B:129:MET:O	1.97	0.47
1:E:184:VAL:HG13	1:E:185:ALA:N	2.29	0.47
1:D:242:GLY:HA2	1:D:275:ARG:NH1	2.29	0.47
1:A:76:LEU:HD21	1:A:138:LEU:HB2	1.96	0.47
1:E:165:TRP:N	1:E:165:TRP:CD1	2.81	0.47
1:E:74:VAL:HG23	1:E:234:LYS:HG2	1.96	0.47
1:A:283:GLY:O	1:A:287:ILE:HG12	2.14	0.47
1:C:62:PRO:O	1:C:184:VAL:HG12	2.15	0.47
1:B:109:THR:C	1:B:111:LYS:H	2.23	0.47
1:C:169:LEU:C	1:C:169:LEU:HD12	2.39	0.47
1:E:256:LYS:O	1:E:256:LYS:HG3	2.15	0.47
1:C:168:GLU:H	1:C:168:GLU:CD	2.22	0.47
1:D:30:LEU:HD11	1:D:58:GLU:OE1	2.15	0.47
1:E:61:ARG:N	1:E:62:PRO:HD3	2.31	0.47
1:B:32:LYS:HE2	1:B:281:GLU:HB3	1.97	0.46
1:B:49:LEU:HD23	1:B:56:ILE:HA	1.97	0.46
1:A:26:HIS:CG	1:A:289:HIS:HB3	2.50	0.46
1:B:261:VAL:HG21	1:B:286:LEU:HD21	1.98	0.46
1:C:250:LEU:HD23	1:C:250:LEU:N	2.29	0.46
1:D:137:LEU:O	1:D:141:THR:HG23	2.15	0.46
1:E:95:ILE:HD12	1:E:122:LEU:HD12	1.97	0.46
1:A:61:ARG:N	1:A:62:PRO:HD3	2.31	0.46
1:D:217:ALA:O	1:D:218:GLY:C	2.58	0.46
1:A:69:MET:HE1	1:A:244:ARG:C	2.41	0.46
1:A:93:ARG:O	1:A:118:THR:HA	2.16	0.46
1:B:32:LYS:HA	1:B:32:LYS:HD2	1.75	0.46
1:B:167:PRO:HD2	1:B:168:GLU:OE2	2.14	0.46
1:E:124:SER:O	1:E:128:THR:OG1	2.27	0.46
1:B:259:ARG:HD2	1:B:290:TRP:CH2	2.51	0.46
1:C:167:PRO:C	1:C:169:LEU:H	2.24	0.46
1:D:225:LEU:HA	1:D:226:PRO:HD3	1.77	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:273:ASP:HB3	1:E:277[B]:ARG:HH21	1.81	0.46
1:C:48:GLU:HB3	1:C:58:GLU:HB2	1.98	0.46
1:C:167:PRO:O	1:C:168:GLU:C	2.57	0.46
1:D:48:GLU:HG2	1:D:57:LEU:HD12	1.98	0.46
1:B:218:GLY:N	1:B:219:PRO:CD	2.79	0.45
1:A:152:LEU:HD13	1:A:160:THR:CG2	2.46	0.45
1:E:47:ILE:HD13	1:E:184:VAL:HA	1.97	0.45
1:C:194:LEU:HD22	1:C:208:ILE:CG2	2.40	0.45
1:A:242:GLY:HA2	1:A:275:ARG:CZ	2.46	0.45
1:C:127:ILE:HA	1:C:234:LYS:HE2	1.97	0.45
1:B:273:ASP:H	1:B:276:ASN:H	1.65	0.45
1:C:99[B]:GLN:H	1:C:99[B]:GLN:HG3	1.24	0.45
1:C:218:GLY:N	1:C:219:PRO:CD	2.80	0.45
1:D:32:LYS:HE2	1:D:281:GLU:CB	2.46	0.45
1:D:216:VAL:HB	3:D:301:MES:H62	1.98	0.45
1:A:36:ALA:O	1:A:40:LEU:HG	2.16	0.45
1:B:171:GLU:HG2	1:B:173:ILE:HG13	1.98	0.45
1:C:67:PRO:O	1:C:69:MET:HE3	2.17	0.45
1:B:26:HIS:CG	1:B:27:PRO:HD2	2.51	0.45
1:C:240:GLU:HB3	5:C:404:HOH:O	2.16	0.45
1:B:148:LEU:HD23	1:B:162:LEU:HD22	1.98	0.45
1:C:26:HIS:CE1	1:C:28:GLU:HG2	2.52	0.45
1:E:89:GLU:CD	1:E:93:ARG:HH21	2.25	0.45
1:B:250:LEU:HD23	1:B:250:LEU:H	1.82	0.44
1:E:70:SER:HB2	1:E:236:GLY:HA2	1.99	0.44
1:A:240:GLU:O	1:A:241:ARG:HB2	2.17	0.44
1:D:51:LEU:HD22	1:D:257:PRO:O	2.16	0.44
1:E:252:PRO:O	1:E:254:ASP:CB	2.27	0.44
1:C:149:THR:O	1:C:153:HIS:HB2	2.17	0.44
1:E:122:LEU:HD22	1:E:134:ALA:HA	1.98	0.44
1:A:26:HIS:N	1:A:58:GLU:HG3	2.33	0.44
1:B:47:ILE:HD13	1:B:184:VAL:HA	2.00	0.44
1:C:252:PRO:O	1:C:254:ASP:CB	2.46	0.44
1:A:199:LEU:HB2	1:A:204:ARG:HG3	1.99	0.44
1:E:80:VAL:O	1:E:84:VAL:HG23	2.17	0.44
1:B:105[B]:TYR:CE1	3:B:301:MES:H51	2.53	0.44
1:D:166:GLU:HA	1:D:167:PRO:HA	1.73	0.44
1:E:26:HIS:HA	1:E:27:PRO:HD3	1.87	0.44
1:B:182:THR:O	1:B:183:PRO:C	2.61	0.44
1:C:137:LEU:HD23	5:C:401:HOH:O	2.18	0.44
1:E:30:LEU:HD12	1:E:30:LEU:C	2.40	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:226:PRO:HG2	1:E:229:TRP:CD1	2.53	0.44
1:C:193:LEU:O	1:C:204:ARG:HG2	2.17	0.44
1:D:244:ARG:HB3	1:D:265:MET:HE3	2.00	0.43
1:C:118:THR:OG1	1:C:121:GLU:HB2	2.17	0.43
1:B:174:PRO:HD3	1:B:241:ARG:HD3	1.99	0.43
1:C:94:ARG:NH1	1:C:115:ASP:O	2.51	0.43
1:E:161:ARG:NH2	1:E:163:ASP:OD2	2.50	0.43
1:A:247:ILE:O	1:A:247:ILE:HG13	2.17	0.43
1:C:61:ARG:N	1:C:62:PRO:CD	2.81	0.43
1:C:166:GLU:HA	1:C:167:PRO:HA	1.75	0.43
1:D:83:ARG:NH1	5:D:407:HOH:O	2.49	0.43
1:B:177:GLU:HA	1:B:180:THR:HG21	2.00	0.43
1:E:208:ILE:O	1:E:209:ASP:C	2.61	0.43
1:D:94:ARG:HD3	1:D:115:ASP:O	2.19	0.43
1:E:51:LEU:H	1:E:259:ARG:HA	1.84	0.43
1:E:81:LEU:O	1:E:84:VAL:HB	2.18	0.43
1:B:29:THR:O	1:B:33:VAL:HG23	2.19	0.43
1:D:233:ASP:HB3	1:D:248:ALA:HB2	2.01	0.43
1:E:217:ALA:HB3	1:E:222:ARG:HE	1.84	0.43
1:C:111:LYS:HD2	5:C:406:HOH:O	2.19	0.43
1:D:26:HIS:CG	1:D:289:HIS:HB3	2.53	0.43
1:D:68:MET:O	1:D:71:THR:OG1	2.19	0.43
1:D:261:VAL:HG21	1:D:286:LEU:CD2	2.49	0.43
1:E:74:VAL:HG21	1:E:247:ILE:HD11	1.99	0.43
1:B:68:MET:O	1:B:69:MET:HB2	2.19	0.42
1:C:174:PRO:O	1:C:175:ASN:HB2	2.19	0.42
1:C:182:THR:OG1	1:C:185:ALA:HB3	2.18	0.42
1:E:236:GLY:O	1:E:244:ARG:HD2	2.20	0.42
1:C:173:ILE:O	1:C:176:ASP:HB3	2.19	0.42
1:A:32:LYS:HE2	1:A:281:GLU:HB3	2.01	0.42
1:A:69:MET:O	1:A:72:PHE:HD1	2.03	0.42
1:C:49:LEU:HD12	1:C:55:LYS:O	2.19	0.42
1:D:206:GLN:O	1:D:209:ASP:HB2	2.20	0.42
1:C:238:GLY:HA3	1:C:243:SER:HB2	2.00	0.42
1:D:261:VAL:HG21	1:D:286:LEU:HD21	2.02	0.42
1:E:148:LEU:HD23	1:E:162:LEU:HD22	2.01	0.42
1:E:184:VAL:HG13	1:E:185:ALA:H	1.85	0.42
1:C:30:LEU:HD23	1:C:30:LEU:HA	1.92	0.42
1:D:48:GLU:HB3	1:D:58:GLU:H	1.85	0.42
1:E:83:ARG:HD3	1:E:142:ILE:HG22	2.02	0.42
1:E:153:HIS:CE1	1:E:158:HIS:CD2	3.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:230:PHE:CE1	1:E:257:PRO:HG3	2.54	0.42
1:C:49:LEU:HD12	1:C:55:LYS:C	2.45	0.42
1:C:51:LEU:HD13	1:C:260:ILE:HG13	2.01	0.42
1:E:220:LEU:HD23	1:E:220:LEU:HA	1.89	0.42
1:A:286:LEU:HD12	1:A:290:TRP:HZ3	1.82	0.42
1:C:225:LEU:HD11	1:C:229:TRP:HB3	2.02	0.42
1:B:94:ARG:NH1	1:B:115:ASP:O	2.48	0.42
1:B:166:GLU:OE1	1:B:167:PRO:HA	2.20	0.42
1:E:211:MET:O	1:E:214:ASP:HB2	2.19	0.42
1:E:225:LEU:HD22	1:E:231:ILE:CD1	2.42	0.42
1:A:47:ILE:HD13	1:A:184:VAL:HA	2.02	0.42
1:A:242:GLY:O	1:A:275:ARG:HD3	2.20	0.42
1:B:153:HIS:HD2	1:B:158:HIS:CG	2.37	0.42
1:C:233:ASP:HB3	1:C:248:ALA:HB2	2.01	0.42
1:E:244:ARG:CB	1:E:265:MET:HE3	2.36	0.42
1:D:47:ILE:CD1	1:D:184:VAL:HA	2.50	0.41
1:A:26:HIS:N	1:A:58:GLU:CG	2.84	0.41
1:A:225:LEU:HA	1:A:226:PRO:HD3	1.91	0.41
1:B:61:ARG:H	1:B:62:PRO:HD3	1.82	0.41
1:C:238:GLY:CA	1:C:243:SER:HB2	2.50	0.41
1:D:69:MET:O	1:D:72:PHE:HD1	2.03	0.41
1:A:130:SER:HG	1:A:234:LYS:HZ1	1.60	0.41
1:D:168:GLU:H	1:D:168:GLU:CD	2.27	0.41
1:D:250:LEU:C	1:D:257:PRO:HA	2.45	0.41
1:C:137:LEU:O	1:C:141:THR:HG23	2.20	0.41
1:C:168:GLU:OE1	1:C:168:GLU:N	2.38	0.41
1:D:32:LYS:HG3	1:D:282:ILE:HG12	2.03	0.41
1:D:184:VAL:HG13	1:D:185:ALA:N	2.36	0.41
1:E:273:ASP:O	1:E:277[B]:ARG:HG3	2.21	0.41
1:D:246:ILE:HG23	1:D:246:ILE:O	2.21	0.41
1:C:106:SER:O	1:C:110:GLU:HG2	2.21	0.41
1:E:65:ARG:HD2	1:E:177:GLU:HG2	2.03	0.41
1:E:174:PRO:O	1:E:175:ASN:HB2	2.21	0.41
1:A:227:ALA:O	1:D:83:ARG:NH2	2.54	0.41
1:D:150:ALA:O	1:D:153[B]:HIS:HB2	2.21	0.41
1:A:126:ALA:O	1:A:130:SER:HA	2.20	0.41
1:B:225:LEU:HA	1:B:226:PRO:HD3	1.82	0.41
1:A:155:MET:HE3	1:A:155:MET:HB2	1.93	0.41
1:B:91:LEU:HD22	1:B:120:GLY:HA2	2.03	0.41
1:B:97:TYR:CD1	1:B:97:TYR:N	2.89	0.41
1:C:61:ARG:N	1:C:62:PRO:HD3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:ILE:O	1:D:231:ILE:HG23	2.19	0.41
1:E:62:PRO:O	1:E:182:THR:HB	2.21	0.41
1:E:233:ASP:HB3	1:E:248:ALA:HB2	2.03	0.41
1:A:168[A]:GLU:H	1:A:168[A]:GLU:HG2	1.55	0.41
1:C:33:VAL:HG22	1:C:282:ILE:HD13	2.01	0.41
1:E:218:GLY:HA3	1:E:219:PRO:HD3	1.89	0.41
1:C:62:PRO:O	1:C:183:PRO:HD2	2.21	0.40
1:D:91:LEU:HA	1:D:119:VAL:HB	2.03	0.40
1:E:30:LEU:CB	1:E:58:GLU:OE1	2.64	0.40
1:E:128:THR:HG22	1:E:210:TRP:O	2.20	0.40
1:E:283:GLY:O	1:E:287:ILE:HG12	2.20	0.40
1:D:146:LYS:HE3	1:D:146:LYS:HB2	1.90	0.40
1:E:155:MET:HE3	1:E:155:MET:HB2	1.92	0.40
1:E:219:PRO:HG2	1:E:276:ASN:HD22	1.85	0.40
1:E:182:THR:O	1:E:183:PRO:C	2.63	0.40
1:E:254:ASP:C	1:E:256:LYS:N	2.77	0.40
1:B:250:LEU:HD23	1:B:250:LEU:N	2.37	0.40
1:C:207:LEU:O	1:C:211:MET:HG3	2.22	0.40
1:D:265:MET:HE2	1:D:279:ILE:HD11	2.02	0.40
1:A:49:LEU:HD12	1:A:56:ILE:HA	2.03	0.40
1:A:90:GLN:HE21	1:A:90:GLN:HB2	1.69	0.40
1:A:174:PRO:HD3	1:A:241:ARG:HD3	2.02	0.40
1:D:205:GLN:O	1:D:208:ILE:HG22	2.22	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:34[A]:LYS:CE	1:E:38:ASP:OD2[1_455]	1.96	0.24
1:E:99:GLN:OE1	1:E:158:HIS:NE2[1_455]	2.05	0.15

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	263/263 (100%)	252 (96%)	10 (4%)	1 (0%)	30	51
1	B	263/263 (100%)	247 (94%)	15 (6%)	1 (0%)	30	51
1	C	264/263 (100%)	251 (95%)	11 (4%)	2 (1%)	16	34
1	D	263/263 (100%)	249 (95%)	13 (5%)	1 (0%)	30	51
1	E	262/263 (100%)	235 (90%)	17 (6%)	10 (4%)	2	5
All	All	1315/1315 (100%)	1234 (94%)	66 (5%)	15 (1%)	11	26

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	226	PRO
1	E	227	ALA
1	E	255	GLY
1	C	168	GLU
1	C	254	ASP
1	E	154	ASN
1	E	218	GLY
1	E	228	GLY
1	E	229	TRP
1	D	105	TYR
1	E	58	GLU
1	E	254	ASP
1	A	105	TYR
1	E	52	ALA
1	B	103	VAL

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	217/215 (101%)	208 (96%)	9 (4%)	27	52
1	B	217/215 (101%)	211 (97%)	6 (3%)	38	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	218/215 (101%)	215 (99%)	3 (1%)	59	80
1	D	217/215 (101%)	214 (99%)	3 (1%)	59	80
1	E	216/215 (100%)	211 (98%)	5 (2%)	44	71
All	All	1085/1075 (101%)	1059 (98%)	26 (2%)	45	70

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	115	ASP
1	A	130	SER
1	A	153[A]	HIS
1	A	153[B]	HIS
1	A	168[A]	GLU
1	A	168[B]	GLU
1	A	171	GLU
1	A	286	LEU
1	B	99	GLN
1	B	114	THR
1	B	123	CYS
1	B	230	PHE
1	B	256	LYS
1	B	273	ASP
1	C	69	MET
1	C	171	GLU
1	C	198	LEU
1	D	39	GLN
1	D	115	ASP
1	D	154	ASN
1	E	28	GLU
1	E	31	VAL
1	E	115	ASP
1	E	123	CYS
1	E	188	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	90	GLN

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Mol	Chain	Res	Type
1	A	206	GLN
1	A	289	HIS
1	B	39	GLN
1	B	153	HIS
1	B	158	HIS
1	B	206	GLN
1	B	269	GLN
1	B	289	HIS
1	C	88	GLN
1	C	158	HIS
1	C	205	GLN
1	C	206	GLN
1	D	39	GLN
1	D	90	GLN
1	D	154	ASN
1	D	158	HIS
1	D	206	GLN
1	E	26	HIS
1	E	39	GLN
1	E	99	GLN
1	E	112	HIS
1	E	158	HIS
1	E	276	ASN
1	E	278	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
3	MES	D	301	-	12,12,12	2.33	1 (8%)	15,16,16	2.48	5 (33%)
4	SO4	C	302	-	4,4,4	0.22	0	6,6,6	0.09	0
3	MES	B	301	-	12,12,12	2.31	1 (8%)	15,16,16	2.45	7 (46%)
3	MES	A	302	-	12,12,12	2.32	1 (8%)	15,16,16	2.47	6 (40%)

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
3	MES	D	301	-	-	5/6/14/14	0/1/1/1
3	MES	B	301	-	-	5/6/14/14	0/1/1/1
3	MES	A	302	-	-	4/6/14/14	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	301	MES	C8-S	-7.79	1.66	1.77
3	A	302	MES	C8-S	-7.76	1.66	1.77
3	B	301	MES	C8-S	-7.73	1.66	1.77

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	301	MES	C5-N4-C3	5.23	120.11	108.84
3	A	302	MES	C5-N4-C3	5.19	120.01	108.84
3	D	301	MES	C5-N4-C3	4.78	119.13	108.84
3	D	301	MES	C2-C3-N4	-4.09	103.90	110.12
3	A	302	MES	C6-C5-N4	-3.60	104.65	110.12
3	D	301	MES	C7-N4-C5	3.51	120.60	111.24
3	B	301	MES	C2-C3-N4	-3.41	104.94	110.12
3	D	301	MES	C7-N4-C3	3.39	120.27	111.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	MES	C7-N4-C5	3.30	120.05	111.24
3	B	301	MES	C7-N4-C5	3.29	120.00	111.24
3	A	302	MES	C7-N4-C3	3.27	119.94	111.24
3	B	301	MES	C7-N4-C3	3.25	119.89	111.24
3	D	301	MES	C6-C5-N4	-3.24	105.20	110.12
3	A	302	MES	C2-C3-N4	-3.21	105.25	110.12
3	B	301	MES	C6-C5-N4	-3.08	105.44	110.12
3	B	301	MES	O2S-S-C8	2.06	109.84	106.73
3	B	301	MES	O1S-S-C8	2.02	109.78	106.73
3	A	302	MES	O1S-S-C8	2.00	109.75	106.73

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	MES	C8-C7-N4-C3
3	D	301	MES	C8-C7-N4-C5
3	D	301	MES	C7-C8-S-O1S
3	D	301	MES	C7-C8-S-O2S
3	D	301	MES	C7-C8-S-O3S
3	B	301	MES	C8-C7-N4-C3
3	B	301	MES	C8-C7-N4-C5
3	A	302	MES	C7-C8-S-O2S
3	B	301	MES	C7-C8-S-O2S
3	A	302	MES	N4-C7-C8-S
3	B	301	MES	N4-C7-C8-S
3	D	301	MES	N4-C7-C8-S
3	B	301	MES	C7-C8-S-O3S
3	A	302	MES	C7-C8-S-O3S

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	301	MES	2	0
4	C	302	SO4	1	0
3	B	301	MES	2	0

5.7 Other polymers [i](#)

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	263/263 (100%)	-0.19	2 (0%) 82 81	10, 20, 34, 44	2 (0%)
1	B	263/263 (100%)	-0.17	2 (0%) 82 81	9, 19, 31, 40	2 (0%)
1	C	263/263 (100%)	0.31	6 (2%) 61 57	18, 35, 55, 68	3 (1%)
1	D	263/263 (100%)	0.46	8 (3%) 52 48	20, 37, 62, 73	2 (0%)
1	E	263/263 (100%)	1.07	32 (12%) 8 7	33, 49, 74, 88	1 (0%)
All	All	1315/1315 (100%)	0.30	50 (3%) 44 39	9, 30, 63, 88	10 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	227	ALA	4.7
1	E	228	GLY	4.1
1	E	105	TYR	3.3
1	D	201	ALA	3.2
1	C	197	GLU	3.2
1	E	221	LEU	3.1
1	E	277[A]	ARG	3.0
1	E	203	SER	3.0
1	D	34	LYS	2.9
1	D	42	GLY	2.9
1	E	26	HIS	2.9
1	E	213	ALA	2.8
1	E	102	LEU	2.7
1	E	27	PRO	2.7
1	E	282	ILE	2.7
1	E	287	ILE	2.7
1	E	290	TRP	2.6
1	D	198	LEU	2.6
1	E	232	ALA	2.6
1	D	281	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	99	GLN	2.5
1	E	193	LEU	2.5
1	C	59	SER	2.5
1	D	153[A]	HIS	2.5
1	D	228	GLY	2.4
1	E	201	ALA	2.4
1	E	56	ILE	2.4
1	C	26	HIS	2.4
1	E	115	ASP	2.4
1	E	289	HIS	2.4
1	E	52	ALA	2.4
1	B	105[A]	TYR	2.3
1	E	194	LEU	2.3
1	E	188	THR	2.3
1	E	223	SER	2.3
1	E	218	GLY	2.3
1	C	52	ALA	2.2
1	E	96	HIS	2.2
1	E	256	LYS	2.2
1	E	103	VAL	2.2
1	B	104[A]	GLU	2.2
1	E	46	TYR	2.1
1	A	153[A]	HIS	2.1
1	E	215	LYS	2.1
1	C	224	ALA	2.1
1	E	217	ALA	2.1
1	E	195	THR	2.1
1	D	277	ARG	2.0
1	E	118	THR	2.0
1	C	277	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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
3	MES	B	301	12/12	0.85	0.23	51,60,66,66	0
4	SO4	C	302	5/5	0.85	0.17	55,55,56,56	5
2	CA	C	301	1/1	0.87	0.08	66,66,66,66	0
3	MES	A	302	12/12	0.88	0.22	52,59,60,61	12
3	MES	D	301	12/12	0.90	0.22	56,63,66,66	12
2	CA	A	301	1/1	0.98	0.04	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.