



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:21 AM UTC

PDB ID : 6IV4 / pdb_00006iv4
Title : Crystal structure of a bacterial Bestrophin homolog from *Klebsiella pneumoniae* with a mutation W252F
Authors : Kittredge, A.; Chen, S.; Yang, T.
Deposited on : 2018-12-02
Resolution : 3.14 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

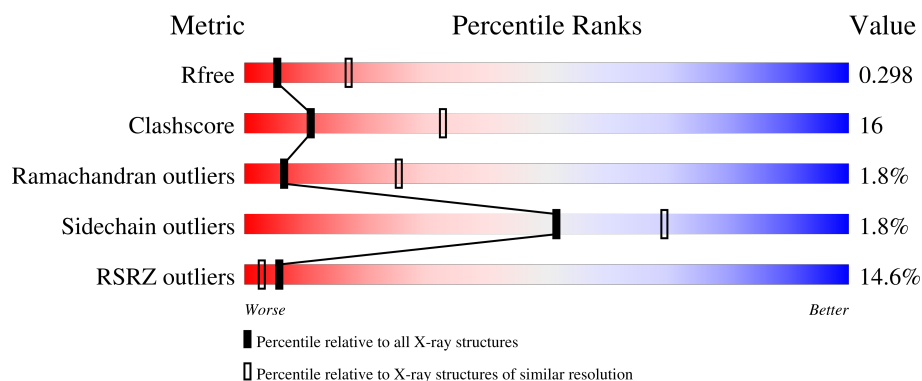
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	296	14% 65% 23% 10%
1	B	296	8% 60% 26% 11%
1	C	296	12% 60% 28% 10%
1	D	296	16% 54% 33% 10%
1	E	296	16% 61% 27% 10%

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	ZN	C	303	-	-	-	X
2	ZN	D	301	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10504 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 Bestrophin homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2108	1366	356	377	9			
1	B	263	Total	C	N	O	S	0	0	0
			2094	1359	353	374	8			
1	C	267	Total	C	N	O	S	0	0	0
			2080	1344	353	374	9			
1	D	267	Total	C	N	O	S	0	0	0
			2099	1359	354	377	9			
1	E	267	Total	C	N	O	S	0	0	0
			2110	1366	358	377	9			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASN	-	expression tag	UNP W1ELP7
A	0	ALA	-	expression tag	UNP W1ELP7
A	252	PHE	TRP	engineered mutation	UNP W1ELP7
B	-1	ASN	-	expression tag	UNP W1ELP7
B	0	ALA	-	expression tag	UNP W1ELP7
B	252	PHE	TRP	engineered mutation	UNP W1ELP7
C	-1	ASN	-	expression tag	UNP W1ELP7
C	0	ALA	-	expression tag	UNP W1ELP7
C	252	PHE	TRP	engineered mutation	UNP W1ELP7
D	-1	ASN	-	expression tag	UNP W1ELP7
D	0	ALA	-	expression tag	UNP W1ELP7
D	252	PHE	TRP	engineered mutation	UNP W1ELP7
E	-1	ASN	-	expression tag	UNP W1ELP7
E	0	ALA	-	expression tag	UNP W1ELP7
E	252	PHE	TRP	engineered mutation	UNP W1ELP7

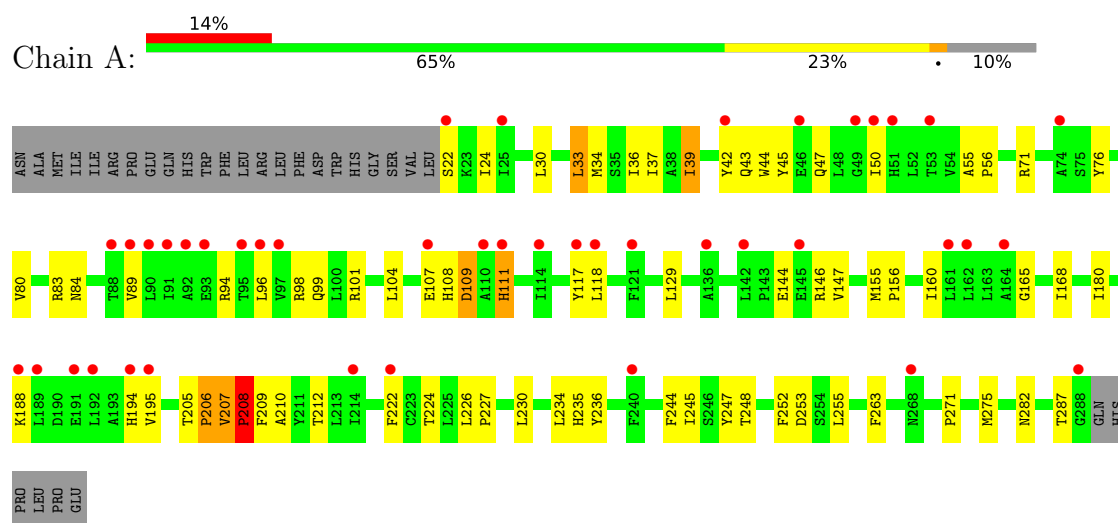
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total 3	Zn 3	0	0
2	B	2	Total 2	Zn 2	0	0
2	C	3	Total 3	Zn 3	0	0
2	D	3	Total 3	Zn 3	0	0
2	E	2	Total 2	Zn 2	0	0

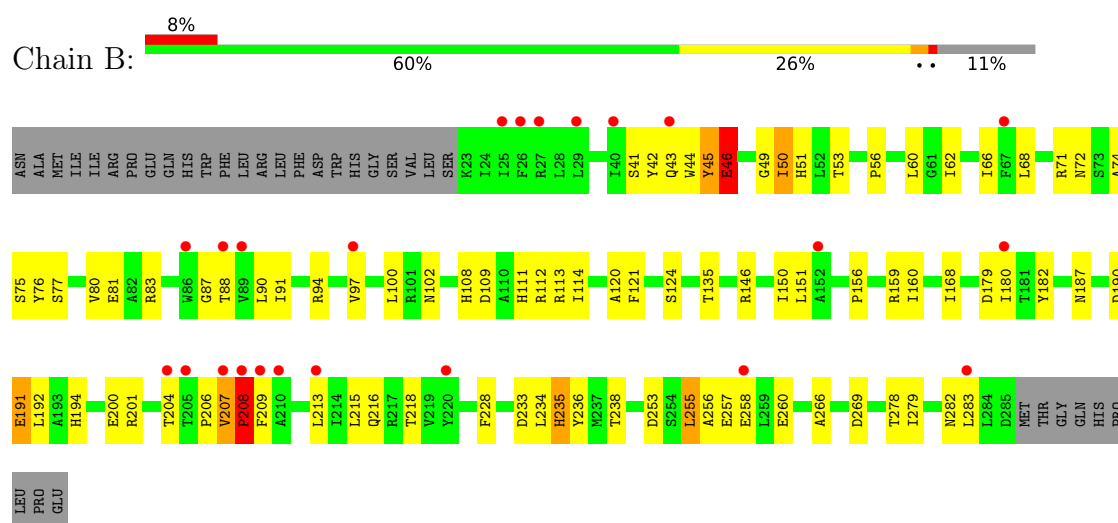
3 Residue-property plots [i](#)

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: Bestrophin homolog

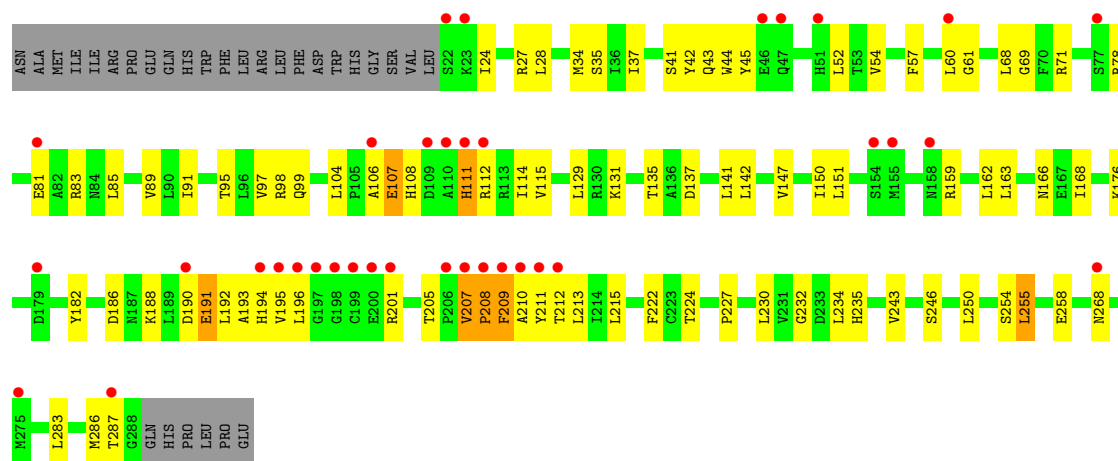


• Molecule 1: Bestrophin homolog

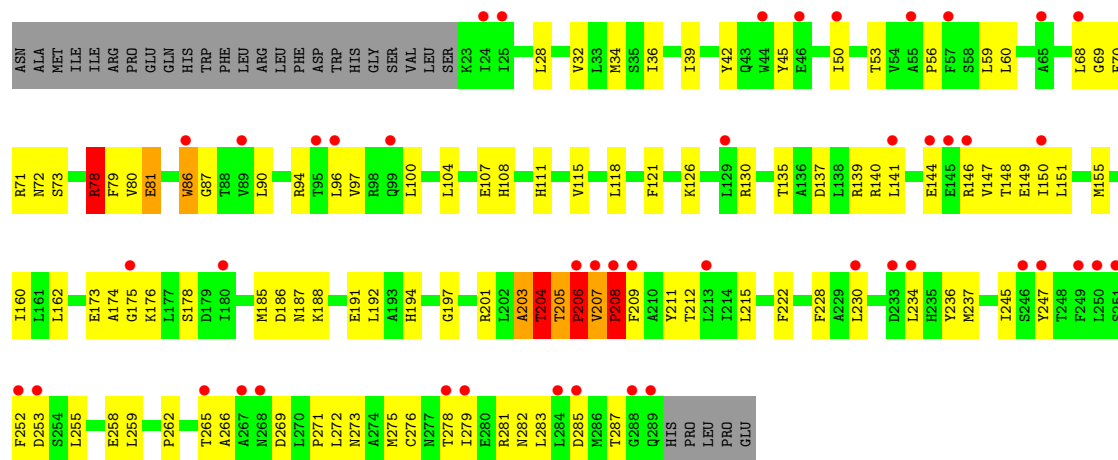


• Molecule 1: Bestrophin homolog

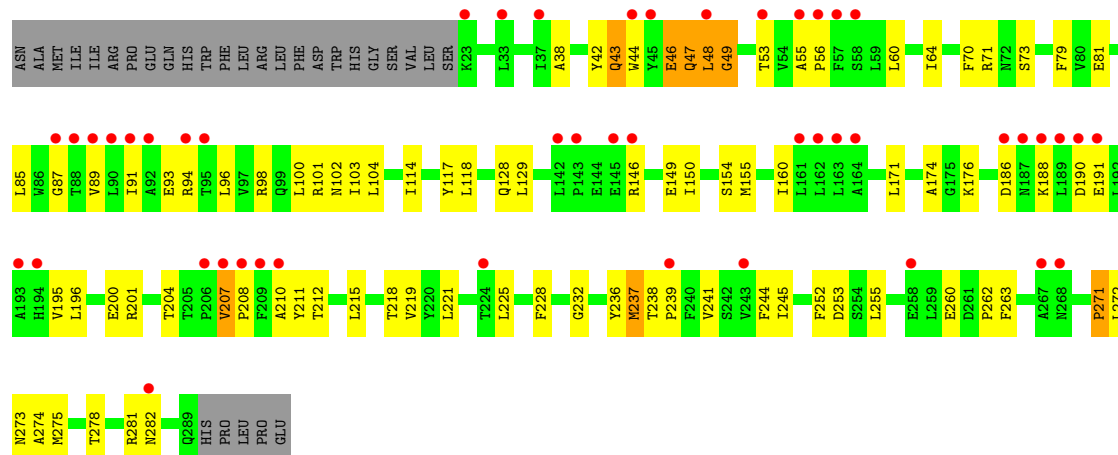




• Molecule 1: Bestrophin homolog



• Molecule 1: Bestrophin homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.01Å 154.19Å 161.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.93 – 3.14 61.93 – 3.14	Depositor EDS
% Data completeness (in resolution range)	98.6 (61.93-3.14) 98.6 (61.93-3.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 3.13Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.237 , 0.295 0.252 , 0.298	Depositor DCC
R_{free} test set	4562 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	104.2	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 25.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	10504	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: 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	A	0.61	1/2153 (0.0%)	0.84	2/2931 (0.1%)
1	B	0.57	1/2139 (0.0%)	0.88	7/2912 (0.2%)
1	C	0.48	1/2121 (0.0%)	0.75	3/2888 (0.1%)
1	D	0.57	2/2143 (0.1%)	0.83	6/2919 (0.2%)
1	E	0.57	3/2154 (0.1%)	0.79	1/2932 (0.0%)
All	All	0.56	8/10710 (0.1%)	0.82	19/14582 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	PRO	C-O	-6.76	1.15	1.23
1	E	191	GLU	CG-CD	6.45	1.68	1.52
1	E	191	GLU	CB-CG	6.22	1.71	1.52
1	D	86	TRP	C-O	-6.11	1.17	1.24
1	C	191	GLU	CG-CD	5.64	1.66	1.52
1	D	81	GLU	C-O	-5.63	1.17	1.24
1	E	271	PRO	C-O	-5.60	1.17	1.24
1	B	191	GLU	CG-CD	5.21	1.65	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	206	PRO	N-CA-C	-10.61	90.61	112.47
1	B	258	GLU	CB-CG-CD	9.20	128.24	112.60
1	D	203	ALA	O-C-N	9.02	134.24	122.42
1	C	235	HIS	CB-CA-C	-8.23	107.09	116.63
1	B	46	GLU	CB-CG-CD	8.18	126.51	112.60
1	A	208	PRO	N-CA-CB	-7.69	95.18	103.25
1	C	111	HIS	CB-CA-C	-7.64	98.89	110.88
1	D	78	ARG	CB-CA-C	-6.75	100.28	110.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	GLU	CB-CA-C	6.50	121.90	110.85
1	E	46	GLU	N-CA-C	-6.03	104.77	111.71
1	C	106	ALA	N-CA-C	5.99	117.61	111.14
1	B	45	TYR	CA-C-O	-5.62	114.92	120.82
1	B	207	VAL	CA-C-N	5.54	126.76	119.84
1	B	207	VAL	C-N-CA	5.54	126.76	119.84
1	D	203	ALA	CA-C-N	5.33	131.72	121.54
1	D	203	ALA	C-N-CA	5.33	131.72	121.54
1	A	109	ASP	N-CA-C	-5.17	105.54	111.07
1	B	255	LEU	N-CA-CB	-5.06	102.69	110.12
1	D	265	THR	CA-CB-OG1	-5.05	102.03	109.60

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	A	2108	0	2142	74	0
1	B	2094	0	2137	89	0
1	C	2080	0	2108	68	0
1	D	2099	0	2125	97	0
1	E	2110	0	2148	73	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	E	2	0	0	0	0
All	All	10504	0	10660	349	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (349) 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:B:46:GLU:OE2	1:B:51:HIS:CE1	1.72	1.41
1:D:201:ARG:CA	1:D:204:THR:OG1	1.65	1.40
1:D:266:ALA:HB3	1:D:269:ASP:OD2	1.20	1.28
1:D:201:ARG:C	1:D:204:THR:OG1	1.86	1.17
1:D:201:ARG:HA	1:D:204:THR:OG1	1.38	1.10
1:B:45:TYR:O	1:B:49:GLY:O	1.73	1.07
1:D:266:ALA:CB	1:D:269:ASP:OD2	2.05	1.04
1:E:236:TYR:O	1:E:239:PRO:HD2	1.59	1.03
1:B:206:PRO:HB3	1:B:209:PHE:HD2	1.18	1.02
1:A:101:ARG:CG	1:A:111:HIS:HE2	1.74	0.99
1:B:206:PRO:HB3	1:B:209:PHE:CD2	1.97	0.99
1:D:81:GLU:OE2	1:D:201:ARG:NH2	1.94	0.98
1:B:46:GLU:OE2	1:B:51:HIS:HE1	1.17	0.97
1:B:208:PRO:HG2	1:D:259:LEU:CD2	1.96	0.95
1:A:101:ARG:HG3	1:A:111:HIS:HE2	1.29	0.95
1:C:268:ASN:HA	1:D:208:PRO:HG2	1.47	0.95
1:B:208:PRO:HG2	1:D:259:LEU:HD21	1.49	0.93
1:B:207:VAL:HG12	1:D:79:PHE:HE1	1.42	0.84
1:B:233:ASP:C	1:B:234:LEU:HD23	2.03	0.84
1:C:68:LEU:HD23	1:C:215:LEU:HD13	1.60	0.84
1:B:206:PRO:CB	1:B:209:PHE:HD2	1.92	0.83
1:B:109:ASP:OD1	1:B:112:ARG:NH2	2.12	0.82
1:A:101:ARG:CG	1:A:111:HIS:NE2	2.43	0.81
1:B:46:GLU:OE2	1:B:51:HIS:NE2	2.13	0.81
1:D:201:ARG:HA	1:D:204:THR:CB	2.10	0.81
1:E:236:TYR:O	1:E:238:THR:N	2.14	0.80
1:B:206:PRO:CB	1:B:209:PHE:CD2	2.64	0.79
1:D:68:LEU:HD23	1:D:215:LEU:HD13	1.67	0.77
1:B:71:ARG:NH2	1:B:253:ASP:OD1	2.17	0.77
1:D:276:CYS:HA	1:D:279:ILE:HD12	1.68	0.76
1:A:271:PRO:HD2	1:A:275:MET:HE3	1.68	0.76
1:B:90:LEU:HD13	1:B:279:ILE:HG12	1.68	0.76
1:C:83:ARG:NH1	1:D:204:THR:O	2.15	0.75
1:A:101:ARG:HG3	1:A:111:HIS:NE2	2.02	0.74
1:A:33:LEU:C	1:A:33:LEU:HD23	2.13	0.74
1:A:101:ARG:HG2	1:A:111:HIS:HE2	1.50	0.74
1:D:266:ALA:HB3	1:D:269:ASP:CG	2.12	0.74
1:B:41:SER:O	1:B:45:TYR:HE1	1.72	0.72
1:C:150:ILE:HA	1:C:159:ARG:HG2	1.71	0.72
1:A:101:ARG:HG2	1:A:111:HIS:NE2	2.04	0.72
1:E:236:TYR:C	1:E:238:THR:H	1.98	0.71
1:B:233:ASP:O	1:B:234:LEU:HD23	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:64:ILE:HD11	1:E:218:THR:HG22	1.73	0.71
1:A:43:GLN:HE21	1:A:44:TRP:CD1	2.09	0.70
1:D:271:PRO:HG2	1:D:275:MET:HE3	1.75	0.69
1:E:81:GLU:OE2	1:E:201:ARG:NH2	2.21	0.68
1:D:104:LEU:HB3	1:D:107:GLU:HG3	1.75	0.68
1:E:94:ARG:NH2	1:E:282:ASN:OD1	2.25	0.68
1:B:207:VAL:HG12	1:D:79:PHE:CE1	2.28	0.68
1:B:68:LEU:HD23	1:B:215:LEU:HD13	1.75	0.67
1:D:278:THR:O	1:D:282:ASN:ND2	2.28	0.67
1:B:108:HIS:CE1	1:B:111:HIS:ND1	2.63	0.67
1:D:201:ARG:O	1:D:204:THR:OG1	2.13	0.66
1:B:46:GLU:CD	1:B:51:HIS:CE1	2.68	0.66
1:A:144:GLU:HA	1:A:147:VAL:HB	1.78	0.65
1:E:71:ARG:NH1	1:E:212:THR:OG1	2.29	0.65
1:B:46:GLU:CD	1:B:51:HIS:NE2	2.54	0.65
1:D:97:VAL:HG21	1:D:283:LEU:HD22	1.79	0.65
1:B:208:PRO:HG2	1:D:259:LEU:HD23	1.77	0.65
1:D:203:ALA:C	1:D:205:THR:H	2.03	0.64
1:D:71:ARG:NH2	1:D:253:ASP:OD1	2.31	0.63
1:B:278:THR:O	1:B:282:ASN:ND2	2.29	0.63
1:E:146:ARG:NH1	1:E:149:GLU:OE1	2.31	0.63
1:E:236:TYR:C	1:E:238:THR:N	2.54	0.63
1:D:211:TYR:CZ	1:D:215:LEU:HG	2.32	0.63
1:C:91:ILE:HD11	1:D:197:GLY:HA3	1.81	0.63
1:A:50:ILE:HG21	1:B:236:TYR:HB2	1.81	0.62
1:D:108:HIS:HA	1:D:111:HIS:CG	2.33	0.62
1:D:108:HIS:CD2	1:D:111:HIS:ND1	2.67	0.62
1:D:137:ASP:OD1	1:D:140:ARG:NH2	2.27	0.62
1:C:89:VAL:HG23	1:C:195:VAL:HG11	1.81	0.62
1:C:209:PHE:HA	1:C:212:THR:HB	1.81	0.62
1:C:207:VAL:O	1:C:209:PHE:N	2.29	0.62
1:C:99:GLN:NE2	1:D:186:ASP:OD2	2.33	0.62
1:A:30:LEU:O	1:A:34:MET:HG2	2.00	0.61
1:A:230:LEU:HB3	1:A:234:LEU:HD12	1.82	0.61
1:E:38:ALA:HB2	1:E:228:PHE:HD1	1.65	0.61
1:D:78:ARG:HH11	1:D:78:ARG:CG	2.13	0.61
1:C:91:ILE:HD13	1:D:194:HIS:HA	1.82	0.61
1:A:99:GLN:NE2	1:E:186:ASP:OD2	2.31	0.61
1:E:150:ILE:HD13	1:E:160:ILE:HG12	1.82	0.61
1:E:271:PRO:HB2	1:E:274:ALA:HB3	1.83	0.60
1:A:108:HIS:HA	1:A:111:HIS:HB2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:53:THR:O	1:D:56:PRO:HD2	2.01	0.60
1:D:146:ARG:NH1	1:D:149:GLU:OE1	2.34	0.60
1:E:271:PRO:O	1:E:273:ASN:N	2.34	0.60
1:C:37:ILE:O	1:C:41:SER:OG	2.12	0.60
1:E:117:TYR:OH	1:E:146:ARG:HG2	2.02	0.59
1:A:98:ARG:NH2	1:E:190:ASP:OD1	2.35	0.59
1:A:156:PRO:O	1:A:160:ILE:HG13	2.02	0.59
1:B:109:ASP:O	1:B:113:ARG:HG3	2.02	0.59
1:B:194:HIS:HE1	1:D:191:GLU:CD	2.10	0.59
1:A:71:ARG:HH11	1:A:212:THR:HG23	1.68	0.58
1:B:200:GLU:O	1:B:204:THR:HB	2.03	0.58
1:D:39:ILE:HD13	1:D:236:TYR:HD1	1.67	0.58
1:A:235:HIS:CD2	1:E:49:GLY:HA3	2.39	0.58
1:A:39:ILE:CD1	1:A:236:TYR:HA	2.34	0.57
1:C:104:LEU:HD21	1:C:176:LYS:O	2.03	0.57
1:D:78:ARG:HH11	1:D:78:ARG:HG3	1.70	0.57
1:C:81:GLU:OE2	1:C:201:ARG:NH2	2.36	0.57
1:E:154:SER:HB2	1:E:155:MET:SD	2.45	0.57
1:B:108:HIS:CD2	1:B:111:HIS:CE1	2.93	0.57
1:D:150:ILE:HD13	1:D:160:ILE:HG12	1.87	0.57
1:D:68:LEU:O	1:D:72:ASN:HB2	2.04	0.56
1:B:135:THR:HG23	1:B:151:LEU:HD11	1.88	0.56
1:B:216:GLN:HG3	1:B:253:ASP:OD2	2.06	0.56
1:D:207:VAL:O	1:D:209:PHE:N	2.39	0.56
1:A:45:TYR:C	1:A:47:GLN:H	2.15	0.55
1:B:108:HIS:HA	1:B:111:HIS:CG	2.41	0.55
1:D:135:THR:HG23	1:D:151:LEU:HD11	1.88	0.55
1:A:96:LEU:HD23	1:A:118:LEU:HD11	1.88	0.55
1:A:107:GLU:O	1:A:109:ASP:N	2.39	0.55
1:B:194:HIS:CE1	1:D:191:GLU:OE2	2.59	0.55
1:C:69:GLY:HA3	1:D:70:PHE:CE2	2.42	0.55
1:D:28:LEU:HD11	1:D:247:TYR:HD1	1.72	0.55
1:C:78:ARG:HH12	1:C:205:THR:CB	2.19	0.55
1:B:194:HIS:HE1	1:D:191:GLU:OE2	1.89	0.55
1:B:42:TYR:HA	1:B:45:TYR:CE1	2.42	0.54
1:E:38:ALA:HB2	1:E:228:PHE:CD1	2.41	0.54
1:B:75:SER:HB3	1:B:260:GLU:HG3	1.88	0.54
1:A:83:ARG:HH12	1:E:204:THR:HG1	1.52	0.54
1:A:222:PHE:O	1:A:226:LEU:HB2	2.07	0.54
1:A:76:TYR:O	1:A:80:VAL:HG23	2.07	0.54
1:B:41:SER:O	1:B:45:TYR:CE1	2.59	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:ASN:OD1	1:E:102:ASN:ND2	2.40	0.54
1:A:224:THR:O	1:A:227:PRO:HD2	2.09	0.53
1:A:80:VAL:O	1:A:84:ASN:ND2	2.39	0.53
1:A:235:HIS:HB3	1:E:49:GLY:O	2.08	0.53
1:C:43:GLN:C	1:C:45:TYR:H	2.17	0.53
1:E:129:LEU:HD13	1:E:263:PHE:CD2	2.43	0.53
1:A:117:TYR:OH	1:A:146:ARG:HD3	2.09	0.53
1:C:35:SER:HB3	1:C:227:PRO:HB3	1.90	0.53
1:D:203:ALA:C	1:D:205:THR:N	2.66	0.53
1:E:89:VAL:HG22	1:E:195:VAL:HG11	1.90	0.53
1:E:43:GLN:O	1:E:46:GLU:HG2	2.08	0.53
1:E:44:TRP:CD1	1:E:44:TRP:H	2.26	0.53
1:C:186:ASP:OD1	1:E:98:ARG:NH1	2.30	0.53
1:C:230:LEU:HB3	1:C:234:LEU:HD12	1.91	0.53
1:A:71:ARG:NH1	1:A:212:THR:HG23	2.23	0.53
1:C:147:VAL:O	1:C:151:LEU:HG	2.09	0.53
1:E:44:TRP:O	1:E:47:GLN:HB3	2.09	0.53
1:C:71:ARG:NH1	1:C:212:THR:HG23	2.24	0.52
1:D:187:ASN:HB2	1:D:188:LYS:HE2	1.90	0.52
1:B:120:ALA:O	1:B:124:SER:HB2	2.10	0.52
1:B:209:PHE:CD1	1:B:209:PHE:C	2.88	0.52
1:D:271:PRO:O	1:D:275:MET:HG3	2.09	0.52
1:C:91:ILE:HG21	1:D:194:HIS:ND1	2.24	0.52
1:A:101:ARG:CG	1:A:111:HIS:CE1	2.93	0.52
1:D:39:ILE:HD13	1:D:236:TYR:CD1	2.45	0.52
1:B:60:LEU:HD13	1:D:245:ILE:HG12	1.92	0.52
1:D:281:ARG:O	1:D:285:ASP:HB2	2.10	0.52
1:C:135:THR:HG23	1:C:151:LEU:HD11	1.92	0.51
1:B:234:LEU:HD23	1:B:234:LEU:N	2.23	0.51
1:B:42:TYR:C	1:B:44:TRP:H	2.18	0.51
1:E:93:GLU:OE1	1:E:118:LEU:HB3	2.10	0.51
1:B:108:HIS:ND1	1:B:111:HIS:ND1	2.58	0.51
1:D:104:LEU:HD21	1:D:176:LYS:O	2.11	0.51
1:D:115:VAL:HG21	1:D:287:THR:HG21	1.91	0.51
1:B:108:HIS:NE2	1:B:111:HIS:CE1	2.79	0.51
1:C:159:ARG:O	1:C:163:LEU:HG	2.10	0.51
1:D:139:ARG:HE	1:D:147:VAL:HG11	1.75	0.51
1:B:206:PRO:CB	1:B:209:PHE:CE2	2.94	0.51
1:C:246:SER:O	1:C:250:LEU:HG	2.11	0.50
1:E:104:LEU:HD21	1:E:176:LYS:O	2.10	0.50
1:A:244:PHE:CZ	1:E:221:LEU:HD23	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:VAL:HG21	1:C:283:LEU:HD22	1.93	0.50
1:C:182:TYR:HE2	1:E:103:ILE:HD11	1.77	0.50
1:E:221:LEU:HG	1:E:225:LEU:HD13	1.92	0.50
1:C:205:THR:C	1:C:207:VAL:H	2.20	0.50
1:C:194:HIS:HA	1:E:91:ILE:HD13	1.94	0.49
1:D:78:ARG:CG	1:D:78:ARG:NH1	2.73	0.49
1:B:235:HIS:O	1:B:238:THR:OG1	2.19	0.49
1:D:68:LEU:CD2	1:D:215:LEU:HD13	2.42	0.49
1:D:108:HIS:CD2	1:D:111:HIS:CE1	3.01	0.49
1:D:86:TRP:O	1:D:87:GLY:C	2.50	0.49
1:C:24:ILE:O	1:C:27:ARG:N	2.46	0.49
1:D:34:MET:HE3	1:D:228:PHE:HE2	1.77	0.49
1:E:55:ALA:HB3	1:E:56:PRO:HD3	1.95	0.49
1:A:89:VAL:HG22	1:A:195:VAL:HG11	1.94	0.49
1:E:100:LEU:HD11	1:E:114:ILE:HD13	1.95	0.49
1:A:39:ILE:HD11	1:A:235:HIS:O	2.12	0.49
1:D:100:LEU:HD23	1:D:185:MET:HE1	1.94	0.49
1:B:121:PHE:CE2	1:B:192:LEU:HD22	2.48	0.49
1:A:71:ARG:NH2	1:A:253:ASP:OD1	2.35	0.48
1:A:244:PHE:CE1	1:E:221:LEU:HD23	2.48	0.48
1:C:188:LYS:NZ	1:C:191:GLU:OE2	2.45	0.48
1:C:208:PRO:C	1:C:210:ALA:H	2.21	0.48
1:A:101:ARG:HD2	1:A:111:HIS:HE1	1.78	0.48
1:C:97:VAL:HG12	1:C:286:MET:HE2	1.95	0.48
1:E:104:LEU:CD1	1:E:171:LEU:HD13	2.44	0.48
1:E:241:VAL:O	1:E:245:ILE:HG13	2.13	0.48
1:A:235:HIS:CG	1:E:49:GLY:HA3	2.49	0.48
1:C:162:LEU:HD21	1:E:101:ARG:HH21	1.79	0.48
1:B:146:ARG:HA	1:B:146:ARG:HD3	1.53	0.48
1:D:139:ARG:HG2	1:D:147:VAL:HG21	1.94	0.48
1:D:252:PHE:O	1:D:255:LEU:HB3	2.14	0.48
1:C:286:MET:HE3	1:D:162:LEU:HD13	1.96	0.48
1:E:236:TYR:O	1:E:237:MET:C	2.49	0.48
1:B:266:ALA:HB3	1:B:269:ASP:OD2	2.15	0.47
1:C:168:ILE:HG22	1:C:182:TYR:CD1	2.49	0.47
1:B:187:ASN:O	1:B:190:ASP:HB2	2.13	0.47
1:C:193:ALA:O	1:C:196:LEU:HB3	2.15	0.47
1:B:215:LEU:HD23	1:B:215:LEU:HA	1.67	0.47
1:B:156:PRO:O	1:B:160:ILE:HG13	2.14	0.47
1:D:42:TYR:HA	1:D:45:TYR:CD1	2.49	0.47
1:B:74:ALA:O	1:B:77:SER:HB2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:GLU:O	1:C:85:LEU:HG	2.15	0.47
1:A:209:PHE:HD2	1:A:209:PHE:O	1.98	0.47
1:A:245:ILE:HG12	1:E:60:LEU:HD13	1.95	0.47
1:B:179:ASP:OD2	1:D:178:SER:OG	2.29	0.47
1:A:42:TYR:O	1:A:45:TYR:HB2	2.15	0.47
1:B:87:GLY:O	1:B:91:ILE:HG13	2.14	0.47
1:B:215:LEU:HD23	1:B:218:THR:HB	1.96	0.47
1:C:107:GLU:H	1:C:107:GLU:HG2	1.63	0.47
1:E:87:GLY:O	1:E:91:ILE:HG13	2.14	0.47
1:A:207:VAL:HG12	1:B:83:ARG:NH2	2.30	0.47
1:B:100:LEU:HD11	1:B:114:ILE:HG21	1.97	0.47
1:E:79:PHE:HD1	1:E:262:PRO:HB3	1.80	0.47
1:A:252:PHE:O	1:A:255:LEU:HB2	2.15	0.47
1:C:162:LEU:HD21	1:E:101:ARG:NH2	2.29	0.47
1:C:150:ILE:HG22	1:C:151:LEU:HD23	1.96	0.46
1:D:69:GLY:O	1:D:73:SER:HB2	2.15	0.46
1:B:81:GLU:OE2	1:B:201:ARG:NH1	2.22	0.46
1:C:60:LEU:HD21	1:E:244:PHE:CE2	2.50	0.46
1:B:190:ASP:O	1:B:194:HIS:HD2	1.98	0.46
1:C:115:VAL:HG21	1:C:287:THR:HG21	1.97	0.46
1:A:180:ILE:HD11	1:B:180:ILE:HD12	1.97	0.46
1:C:60:LEU:HD23	1:C:222:PHE:HD1	1.81	0.46
1:C:168:ILE:HG22	1:C:182:TYR:HD1	1.79	0.46
1:D:230:LEU:HB3	1:D:234:LEU:HD12	1.97	0.46
1:A:33:LEU:C	1:A:33:LEU:CD2	2.85	0.46
1:A:39:ILE:HD11	1:A:236:TYR:HA	1.97	0.46
1:E:208:PRO:HG2	1:E:211:TYR:H	1.80	0.46
1:D:139:ARG:HG2	1:D:147:VAL:HG11	1.98	0.46
1:D:209:PHE:O	1:D:212:THR:HG22	2.15	0.46
1:B:45:TYR:OH	1:B:228:PHE:HD1	1.99	0.46
1:C:95:THR:O	1:C:99:GLN:HG3	2.16	0.46
1:C:190:ASP:C	1:C:192:LEU:N	2.71	0.46
1:E:96:LEU:HD23	1:E:118:LEU:HD21	1.98	0.46
1:E:85:LEU:HD23	1:E:85:LEU:HA	1.78	0.45
1:B:75:SER:CB	1:B:260:GLU:HG3	2.45	0.45
1:A:188:LYS:HA	1:A:188:LYS:HD3	1.81	0.45
1:B:45:TYR:C	1:B:49:GLY:O	2.56	0.45
1:B:208:PRO:CG	1:D:259:LEU:HD21	2.34	0.45
1:E:196:LEU:O	1:E:200:GLU:HG3	2.16	0.45
1:C:42:TYR:CZ	1:C:232:GLY:HA2	2.51	0.45
1:A:55:ALA:HB3	1:A:56:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:ARG:NH1	1:E:204:THR:OG1	2.35	0.45
1:A:111:HIS:CE1	1:A:287:THR:HG22	2.52	0.45
1:D:96:LEU:HD23	1:D:118:LEU:HD11	1.99	0.45
1:A:180:ILE:HD13	1:A:180:ILE:HA	1.77	0.45
1:A:236:TYR:HD2	1:E:48:LEU:HB3	1.81	0.45
1:D:173:GLU:C	1:D:175:GLY:H	2.25	0.45
1:E:53:THR:HB	1:E:56:PRO:HD2	1.98	0.45
1:A:94:ARG:NH2	1:A:282:ASN:OD1	2.42	0.44
1:D:155:MET:SD	1:D:155:MET:N	2.91	0.44
1:D:126:LYS:HE3	1:D:273:ASN:ND2	2.32	0.44
1:E:71:ARG:NH2	1:E:253:ASP:OD1	2.50	0.44
1:A:22:SER:C	1:A:24:ILE:H	2.24	0.44
1:B:108:HIS:CG	1:B:111:HIS:ND1	2.86	0.44
1:A:108:HIS:CA	1:A:111:HIS:HB2	2.46	0.44
1:A:129:LEU:HD13	1:A:263:PHE:CD1	2.53	0.44
1:A:89:VAL:CG2	1:A:195:VAL:HG11	2.48	0.44
1:C:129:LEU:C	1:C:131:LYS:H	2.26	0.44
1:B:194:HIS:HE1	1:D:191:GLU:OE1	2.01	0.43
1:C:111:HIS:O	1:C:114:ILE:N	2.46	0.43
1:A:180:ILE:HD12	1:A:180:ILE:HG23	1.73	0.43
1:B:50:ILE:HA	1:D:237:MET:HG2	1.99	0.43
1:C:191:GLU:O	1:C:195:VAL:HG23	2.17	0.43
1:B:207:VAL:O	1:B:207:VAL:HG23	2.17	0.43
1:D:188:LYS:HA	1:D:188:LYS:HD3	1.89	0.43
1:E:104:LEU:HD12	1:E:171:LEU:HD13	2.01	0.43
1:E:252:PHE:O	1:E:255:LEU:HB3	2.18	0.43
1:D:115:VAL:HG21	1:D:287:THR:CG2	2.48	0.43
1:D:144:GLU:O	1:D:148:THR:HG23	2.19	0.43
1:E:212:THR:HG21	1:E:260:GLU:OE1	2.18	0.43
1:A:43:GLN:HE21	1:A:44:TRP:NE1	2.17	0.43
1:A:194:HIS:HA	1:B:91:ILE:HG21	2.00	0.43
1:B:97:VAL:HG21	1:B:283:LEU:HD22	1.99	0.43
1:C:28:LEU:HD22	1:C:243:VAL:HG23	2.00	0.43
1:C:78:ARG:HE	1:C:78:ARG:HB3	1.63	0.43
1:B:72:ASN:OD1	1:B:256:ALA:HB2	2.18	0.43
1:C:205:THR:O	1:C:207:VAL:N	2.47	0.43
1:D:206:PRO:HD2	1:D:206:PRO:O	2.17	0.43
1:C:135:THR:C	1:C:137:ASP:H	2.27	0.43
1:C:52:LEU:HD23	1:C:52:LEU:HA	1.86	0.43
1:B:42:TYR:O	1:B:44:TRP:N	2.52	0.42
1:E:278:THR:HA	1:E:281:ARG:HE	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:GLY:HA3	1:D:59:LEU:HD11	2.00	0.42
1:D:205:THR:HA	1:D:206:PRO:HD3	1.85	0.42
1:B:213:LEU:HD13	1:B:213:LEU:HA	1.84	0.42
1:E:100:LEU:HD23	1:E:100:LEU:HA	1.89	0.42
1:C:34:MET:HE2	1:C:224:THR:HB	2.01	0.42
1:D:108:HIS:CG	1:D:111:HIS:ND1	2.87	0.42
1:D:162:LEU:HD12	1:D:162:LEU:HA	1.90	0.42
1:A:165:GLY:HA3	1:B:102:ASN:OD1	2.19	0.42
1:A:247:TYR:O	1:A:248:THR:C	2.63	0.42
1:B:206:PRO:HG2	1:B:209:PHE:HE2	1.84	0.42
1:B:255:LEU:HD12	1:B:255:LEU:HA	1.74	0.42
1:D:140:ARG:HG2	1:D:141:LEU:HG	2.01	0.42
1:E:129:LEU:HD13	1:E:263:PHE:CE2	2.55	0.42
1:A:50:ILE:CG2	1:B:236:TYR:HB2	2.48	0.42
1:A:155:MET:HE2	1:B:94:ARG:HH22	1.85	0.42
1:B:168:ILE:HG22	1:B:182:TYR:CD1	2.54	0.42
1:E:42:TYR:CE2	1:E:232:GLY:HA2	2.54	0.42
1:B:88:THR:HA	1:B:91:ILE:HD12	2.00	0.42
1:C:54:VAL:O	1:C:57:PHE:N	2.52	0.42
1:D:111:HIS:N	1:D:111:HIS:CD2	2.88	0.42
1:E:271:PRO:HD2	1:E:275:MET:HE3	2.01	0.42
1:B:150:ILE:HA	1:B:159:ARG:HG2	2.01	0.42
1:D:90:LEU:HG	1:D:94:ARG:NH1	2.35	0.42
1:E:215:LEU:O	1:E:219:VAL:HG23	2.19	0.42
1:A:155:MET:HE2	1:B:94:ARG:NH2	2.34	0.42
1:C:211:TYR:O	1:C:215:LEU:N	2.43	0.42
1:B:53:THR:O	1:B:56:PRO:HD2	2.20	0.41
1:D:203:ALA:O	1:D:205:THR:N	2.38	0.41
1:A:76:TYR:CE2	1:A:80:VAL:HG21	2.56	0.41
1:C:255:LEU:HD12	1:C:255:LEU:HA	1.79	0.41
1:D:130:ARG:NH2	1:D:272:LEU:HB2	2.35	0.41
1:B:108:HIS:HA	1:B:111:HIS:CD2	2.55	0.41
1:C:141:LEU:O	1:C:142:LEU:HD23	2.20	0.41
1:E:93:GLU:CD	1:E:118:LEU:HB3	2.45	0.41
1:E:207:VAL:HB	1:E:208:PRO:HD3	2.01	0.41
1:E:211:TYR:CE2	1:E:215:LEU:HD12	2.55	0.41
1:A:101:ARG:HD2	1:A:111:HIS:CE1	2.55	0.41
1:B:76:TYR:CE2	1:B:80:VAL:HG21	2.55	0.41
1:E:174:ALA:HB3	1:E:176:LYS:HG3	2.02	0.41
1:A:45:TYR:C	1:A:47:GLN:N	2.78	0.41
1:A:83:ARG:NH1	1:E:204:THR:HG1	2.16	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:HIS:HE1	1:B:191:GLU:OE2	2.03	0.41
1:C:107:GLU:CD	1:C:176:LYS:HD3	2.43	0.41
1:D:32:VAL:O	1:D:36:ILE:HG12	2.20	0.41
1:C:254:SER:O	1:C:258:GLU:HG3	2.21	0.41
1:C:85:LEU:HB3	1:C:195:VAL:HG13	2.01	0.41
1:C:98:ARG:HG3	1:D:162:LEU:HD12	2.02	0.41
1:E:79:PHE:CD1	1:E:262:PRO:HB3	2.55	0.41
1:C:213:LEU:HD12	1:C:213:LEU:HA	1.71	0.41
1:D:50:ILE:HG13	1:D:50:ILE:O	2.20	0.41
1:B:194:HIS:CE1	1:D:191:GLU:CD	2.97	0.41
1:B:62:ILE:O	1:B:66:ILE:HG13	2.20	0.40
1:E:70:PHE:O	1:E:73:SER:HB3	2.21	0.40
1:A:104:LEU:HB3	1:A:107:GLU:HG2	2.01	0.40
1:A:168:ILE:HG23	1:A:168:ILE:HD12	1.82	0.40
1:D:108:HIS:HA	1:D:111:HIS:CD2	2.56	0.40
1:D:121:PHE:CE2	1:D:192:LEU:HD22	2.56	0.40
1:D:60:LEU:HD23	1:D:222:PHE:HD1	1.87	0.40
1:E:128:GLN:HE22	1:E:196:LEU:HD11	1.86	0.40
1:E:188:LYS:HA	1:E:188:LYS:HD3	1.89	0.40
1:A:205:THR:N	1:A:206:PRO:CD	2.84	0.40

There are no symmetry-related clashes.

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	265/296 (90%)	241 (91%)	22 (8%)	2 (1%)	16	44
1	B	261/296 (88%)	245 (94%)	13 (5%)	3 (1%)	11	36
1	C	265/296 (90%)	239 (90%)	21 (8%)	5 (2%)	6	24
1	D	265/296 (90%)	238 (90%)	20 (8%)	7 (3%)	4	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	265/296 (90%)	240 (91%)	18 (7%)	7 (3%)	4	18
All	All	1321/1480 (89%)	1203 (91%)	94 (7%)	24 (2%)	6	25

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	210	ALA
1	B	208	PRO
1	D	208	PRO
1	B	50	ILE
1	E	237	MET
1	A	208	PRO
1	B	43	GLN
1	C	209	PHE
1	C	255	LEU
1	E	48	LEU
1	C	44	TRP
1	D	204	THR
1	D	262	PRO
1	E	43	GLN
1	E	49	GLY
1	E	272	LEU
1	C	207	VAL
1	D	174	ALA
1	E	210	ALA
1	D	206	PRO
1	D	258	GLU
1	C	208	PRO
1	E	207	VAL
1	D	205	THR

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	228/259 (88%)	221 (97%)	7 (3%)	35	60
1	B	228/259 (88%)	224 (98%)	4 (2%)	51	69
1	C	222/259 (86%)	219 (99%)	3 (1%)	59	72
1	D	226/259 (87%)	221 (98%)	5 (2%)	45	67
1	E	228/259 (88%)	227 (100%)	1 (0%)	84	84
All	All	1132/1295 (87%)	1112 (98%)	20 (2%)	51	69

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

Mol	Chain	Res	Type
1	A	33	LEU
1	A	36	ILE
1	A	37	ILE
1	A	39	ILE
1	A	111	HIS
1	A	207	VAL
1	A	208	PRO
1	B	46	GLU
1	B	208	PRO
1	B	235	HIS
1	B	257	GLU
1	C	107	GLU
1	C	108	HIS
1	C	112	ARG
1	D	78	ARG
1	D	80	VAL
1	D	204	THR
1	D	207	VAL
1	D	208	PRO
1	E	47	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	84	ASN
1	A	235	HIS
1	B	51	HIS
1	B	102	ASN
1	B	108	HIS

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Mol	Chain	Res	Type
1	B	194	HIS
1	D	31	ASN
1	D	108	HIS
1	D	273	ASN
1	D	289	GLN
1	E	51	HIS
1	E	99	GLN
1	E	289	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 13 ligands modelled in this entry, 13 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 ⓘ

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	267/296 (90%)	0.86	42 (15%) 5 2	68, 90, 119, 160	0
1	B	263/296 (88%)	0.58	23 (8%) 16 8	70, 91, 128, 162	0
1	C	267/296 (90%)	0.93	36 (13%) 7 3	72, 105, 134, 152	0
1	D	267/296 (90%)	0.83	46 (17%) 4 2	74, 99, 130, 158	0
1	E	267/296 (90%)	0.99	47 (17%) 4 2	66, 88, 118, 139	0
All	All	1331/1480 (89%)	0.84	194 (14%) 6 3	66, 95, 130, 162	0

All (194) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	210	ALA	11.1
1	C	109	ASP	10.4
1	E	209	PHE	10.0
1	B	26	PHE	9.8
1	E	88	THR	8.9
1	C	198	GLY	8.7
1	D	145	GLU	8.6
1	E	145	GLU	8.2
1	C	207	VAL	7.9
1	C	208	PRO	7.7
1	C	112	ARG	7.3
1	C	194	HIS	6.9
1	E	187	ASN	6.9
1	C	197	GLY	6.5
1	B	220	TYR	6.2
1	A	92	ALA	6.1
1	E	91	ILE	6.0
1	E	90	LEU	5.8
1	A	49	GLY	5.6
1	C	209	PHE	5.5

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Mol	Chain	Res	Type	RSRZ
1	E	94	ARG	5.5
1	C	199	CYS	5.5
1	B	25	ILE	5.5
1	B	27	ARG	5.4
1	E	268	ASN	5.4
1	E	92	ALA	5.4
1	C	206	PRO	5.4
1	C	111	HIS	5.3
1	A	111	HIS	5.3
1	D	288	GLY	5.2
1	E	258	GLU	5.0
1	C	47	GLN	4.9
1	C	196	LEU	4.8
1	C	200	GLU	4.8
1	B	204	THR	4.7
1	B	213	LEU	4.6
1	B	88	THR	4.6
1	A	191	GLU	4.5
1	E	208	PRO	4.5
1	B	207	VAL	4.5
1	A	51	HIS	4.4
1	D	55	ALA	4.4
1	E	95	THR	4.4
1	E	207	VAL	4.4
1	D	285	ASP	4.3
1	C	155	MET	4.3
1	D	267	ALA	4.3
1	A	93	GLU	4.3
1	E	194	HIS	4.2
1	C	211	TYR	4.2
1	E	89	VAL	4.1
1	D	44	TRP	4.1
1	C	268	ASN	4.1
1	E	191	GLU	4.1
1	A	195	VAL	4.0
1	E	56	PRO	4.0
1	D	86	TRP	3.9
1	C	190	ASP	3.9
1	E	58	SER	3.9
1	A	42	TYR	3.9
1	B	29	LEU	3.8
1	E	163	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	43	GLN	3.7
1	E	161	LEU	3.7
1	C	51	HIS	3.7
1	E	188	LYS	3.7
1	C	77	SER	3.7
1	D	206	PRO	3.6
1	B	205	THR	3.6
1	A	110	ALA	3.5
1	E	55	ALA	3.5
1	E	190	ASP	3.5
1	A	136	ALA	3.3
1	D	234	LEU	3.3
1	D	99	GLN	3.3
1	D	230	LEU	3.3
1	A	117	TYR	3.2
1	A	114	ILE	3.2
1	D	68	LEU	3.2
1	E	48	LEU	3.2
1	C	158	ASN	3.2
1	A	97	VAL	3.2
1	D	289	GLN	3.2
1	C	201	ARG	3.2
1	A	192	LEU	3.1
1	D	207	VAL	3.1
1	D	144	GLU	3.1
1	E	23	LYS	3.1
1	A	95	THR	3.1
1	D	284	LEU	3.0
1	C	46	GLU	3.0
1	D	233	ASP	3.0
1	E	162	LEU	3.0
1	E	57	PHE	2.9
1	D	250	LEU	2.9
1	E	210	ALA	2.9
1	E	87	GLY	2.9
1	C	287	THR	2.9
1	A	89	VAL	2.9
1	B	208	PRO	2.9
1	B	283	LEU	2.9
1	C	212	THR	2.8
1	A	145	GLU	2.8
1	C	60	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	25	ILE	2.8
1	E	193	ALA	2.8
1	D	208	PRO	2.8
1	E	143	PRO	2.8
1	D	95	THR	2.8
1	C	154	SER	2.7
1	A	164	ALA	2.7
1	C	179	ASP	2.7
1	B	40	ILE	2.7
1	E	146	ARG	2.6
1	B	86	TRP	2.6
1	E	45	TYR	2.6
1	C	195	VAL	2.6
1	C	22	SER	2.6
1	A	188	LYS	2.6
1	E	33	LEU	2.6
1	A	162	LEU	2.6
1	A	107	GLU	2.6
1	D	268	ASN	2.5
1	A	88	THR	2.5
1	D	180	ILE	2.5
1	A	222	PHE	2.5
1	D	246	SER	2.5
1	A	46	GLU	2.5
1	D	150	ILE	2.5
1	B	152	ALA	2.5
1	D	251	SER	2.5
1	D	57	PHE	2.5
1	D	278	THR	2.5
1	E	243	VAL	2.5
1	E	224	THR	2.4
1	B	180	ILE	2.4
1	A	121	PHE	2.4
1	D	252	PHE	2.4
1	D	265	THR	2.4
1	D	89	VAL	2.4
1	D	279	ILE	2.4
1	E	164	ALA	2.4
1	D	249	PHE	2.4
1	B	97	VAL	2.4
1	E	282	ASN	2.4
1	B	210	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	106	ALA	2.4
1	C	23	LYS	2.4
1	B	209	PHE	2.4
1	A	268	ASN	2.3
1	E	206	PRO	2.3
1	E	267	ALA	2.3
1	D	247	TYR	2.3
1	D	46	GLU	2.3
1	D	146	ARG	2.3
1	A	194	HIS	2.3
1	D	24	ILE	2.3
1	D	175	GLY	2.3
1	D	253	ASP	2.3
1	A	53	THR	2.2
1	A	50	ILE	2.2
1	A	90	LEU	2.2
1	A	96	LEU	2.2
1	E	189	LEU	2.2
1	C	275	MET	2.2
1	D	213	LEU	2.2
1	B	258	GLU	2.2
1	D	96	LEU	2.2
1	E	186	ASP	2.2
1	D	209	PHE	2.2
1	A	142	LEU	2.1
1	A	161	LEU	2.1
1	E	44	TRP	2.1
1	A	288	GLY	2.1
1	B	89	VAL	2.1
1	A	25	ILE	2.1
1	A	189	LEU	2.1
1	D	50	ILE	2.1
1	E	37	ILE	2.1
1	A	240	PHE	2.1
1	D	141	LEU	2.1
1	E	239	PRO	2.1
1	A	118	LEU	2.1
1	D	129	LEU	2.1
1	E	142	LEU	2.1
1	E	53	THR	2.1
1	A	74	ALA	2.1
1	C	110	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	67	PHE	2.0
1	A	22	SER	2.0
1	D	65	ALA	2.0
1	C	81	GLU	2.0
1	A	91	ILE	2.0
1	A	214	ILE	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 ⓘ

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	ZN	D	301	1/1	0.33	0.69	90,90,90,90	0
2	ZN	C	303	1/1	0.47	0.72	110,110,110,110	0
2	ZN	B	302	1/1	0.50	0.31	79,79,79,79	0
2	ZN	E	301	1/1	0.64	0.23	111,111,111,111	0
2	ZN	D	302	1/1	0.66	0.19	112,112,112,112	0
2	ZN	A	301	1/1	0.67	0.22	82,82,82,82	0
2	ZN	B	301	1/1	0.72	0.14	93,93,93,93	0
2	ZN	C	301	1/1	0.73	0.25	89,89,89,89	0
2	ZN	A	303	1/1	0.74	0.13	95,95,95,95	0
2	ZN	E	302	1/1	0.75	0.16	93,93,93,93	0
2	ZN	A	302	1/1	0.82	0.14	90,90,90,90	0
2	ZN	C	302	1/1	0.85	0.31	84,84,84,84	0
2	ZN	D	303	1/1	0.88	0.11	114,114,114,114	0

6.5 Other polymers ⓘ

There are no such residues in this entry.