



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 09:20 PM UTC

PDB ID : 6IVL / pdb_00006ivl
Title : Crystal structure of a membrane protein L259A
Authors : Kittredge, A.; Fukuda, F.; Zhang, Y.; Yang, T.
Deposited on : 2018-12-04
Resolution : 3.40 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

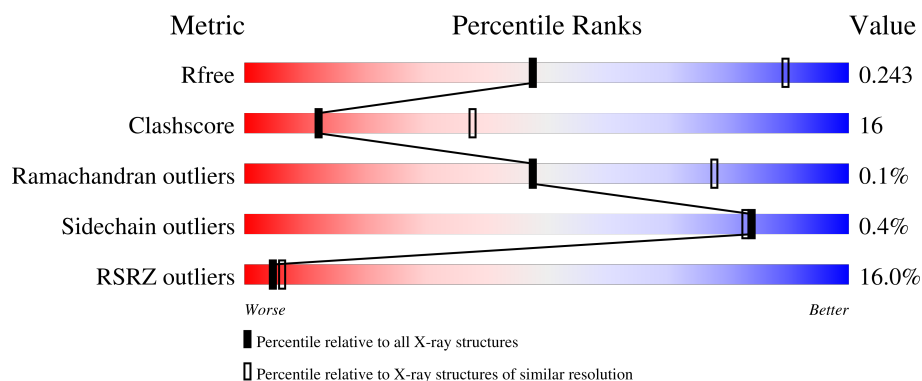
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	297	8% 62% 28% 10%
1	B	297	14% 61% 29% • 9%
1	C	297	12% 63% 28% 9%
1	D	297	24% 68% 22% • 10%
1	E	297	15% 70% 21% 9%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10735 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 Ibestrophin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2132	1380	361	382	9			
1	B	270	Total	C	N	O	S	0	0	0
			2149	1391	365	384	9			
1	C	269	Total	C	N	O	S	0	0	0
			2137	1382	364	382	9			
1	D	268	Total	C	N	O	S	0	0	0
			2132	1380	361	382	9			
1	E	270	Total	C	N	O	S	0	0	0
			2149	1393	364	383	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP W9BH30
A	-1	ASN	-	expression tag	UNP W9BH30
A	0	ALA	-	expression tag	UNP W9BH30
A	259	ALA	LEU	engineered mutation	UNP W9BH30
B	-2	SER	-	expression tag	UNP W9BH30
B	-1	ASN	-	expression tag	UNP W9BH30
B	0	ALA	-	expression tag	UNP W9BH30
B	259	ALA	LEU	engineered mutation	UNP W9BH30
C	-2	SER	-	expression tag	UNP W9BH30
C	-1	ASN	-	expression tag	UNP W9BH30
C	0	ALA	-	expression tag	UNP W9BH30
C	259	ALA	LEU	engineered mutation	UNP W9BH30
D	-2	SER	-	expression tag	UNP W9BH30
D	-1	ASN	-	expression tag	UNP W9BH30
D	0	ALA	-	expression tag	UNP W9BH30
D	259	ALA	LEU	engineered mutation	UNP W9BH30
E	-2	SER	-	expression tag	UNP W9BH30
E	-1	ASN	-	expression tag	UNP W9BH30
E	0	ALA	-	expression tag	UNP W9BH30

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Chain	Residue	Modelled	Actual	Comment	Reference
E	259	ALA	LEU	engineered mutation	UNP W9BH30

- 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	4	Total 4	Zn 4	0	0
2	C	3	Total 3	Zn 3	0	0
2	D	3	Total 3	Zn 3	0	0
2	E	3	Total 3	Zn 3	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Cl 1	0	0
3	C	1	Total 1	Cl 1	0	0
3	D	2	Total 2	Cl 2	0	0
3	E	2	Total 2	Cl 2	0	0

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

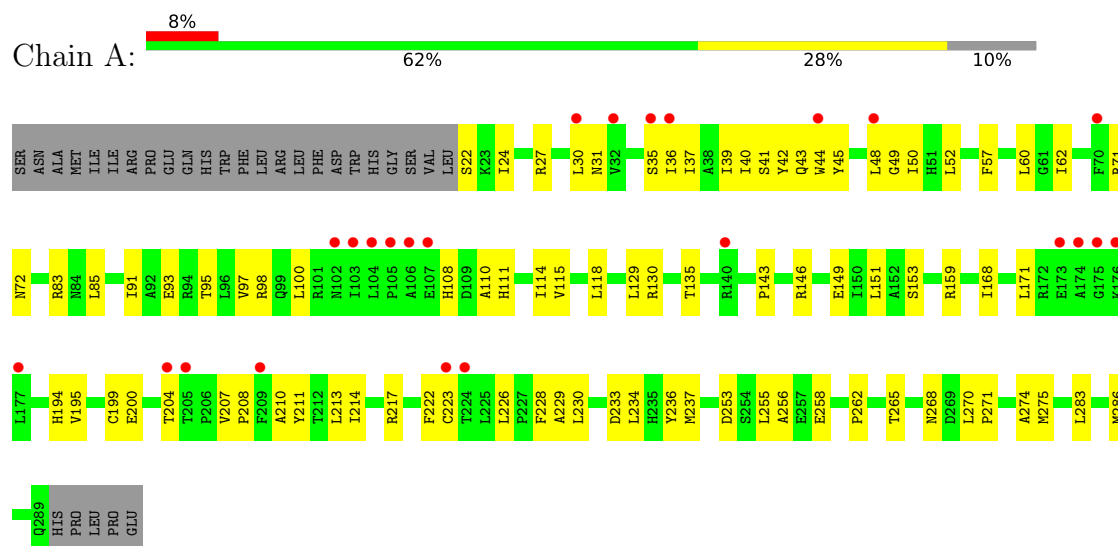
- Molecule 5 is water.

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

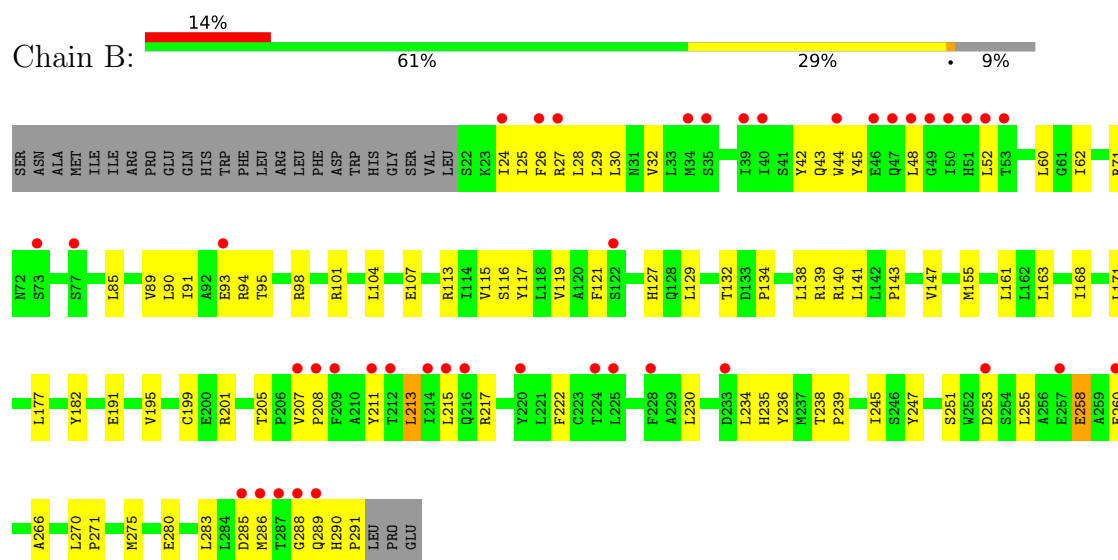
3 Residue-property plots [i](#)

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: Ibestrophin

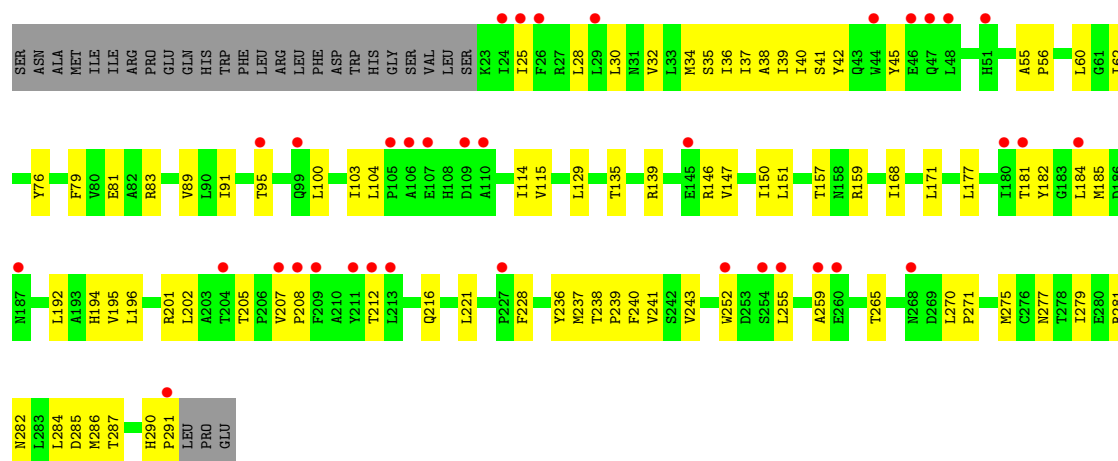


• Molecule 1: Ibestrophin

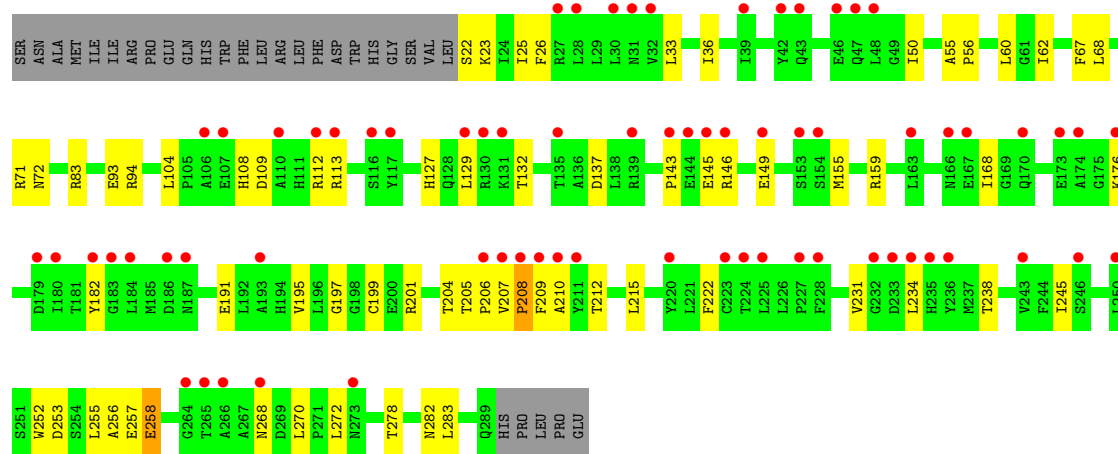


• Molecule 1: Ibestrophin

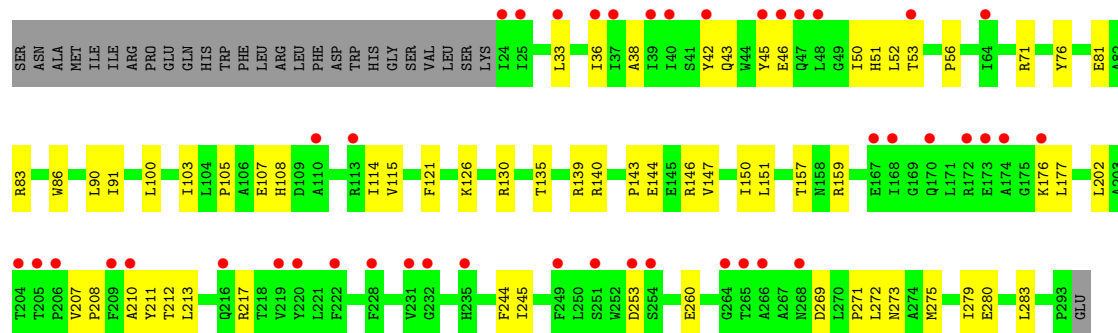




• Molecule 1: Ibestrophin



• Molecule 1: Ibestrophin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.35Å 160.54Å 161.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 3.40 48.80 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.80-3.40) 99.9 (48.80-3.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 3.40Å)	Xtriage
Refinement program	REFMAC 5.8.0238, PHENIX	Depositor
R, R_{free}	0.209 , 0.244 0.208 , 0.243	Depositor DCC
R_{free} test set	2088 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	109.9	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	10735	wwPDB-VP
Average B, all atoms (Å ²)	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: CL, ZN, ACY

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	0.49	0/2178	0.69	0/2964
1	B	0.45	0/2197	0.72	0/2991
1	C	0.47	0/2184	0.64	0/2974
1	D	0.43	0/2178	0.61	0/2964
1	E	0.47	0/2198	0.64	0/2995
All	All	0.46	0/10935	0.66	0/14888

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

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	43	GLN	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2132	0	2175	86	0
1	B	2149	0	2189	103	0
1	C	2137	0	2177	76	0
1	D	2132	0	2175	71	0
1	E	2149	0	2189	60	0
2	A	3	0	0	0	0
2	B	4	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	E	3	0	0	0	0
3	A	1	0	0	1	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	1	0
4	C	4	0	3	0	0
5	A	2	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
All	All	10735	0	10908	336	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 (336) 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:213:LEU:HD11	1:B:217:ARG:CD	1.78	1.14
1:B:213:LEU:HD11	1:B:217:ARG:HD2	1.22	1.10
1:B:213:LEU:CD1	1:B:217:ARG:HD2	1.82	1.07
1:B:211:TYR:CD2	1:D:255:LEU:HD23	1.92	1.04
1:B:213:LEU:CD1	1:B:217:ARG:CD	2.37	1.02
1:C:236:TYR:O	1:C:239:PRO:HD2	1.68	0.93
1:B:44:TRP:O	1:B:48:LEU:CD1	2.18	0.91
1:B:44:TRP:O	1:B:48:LEU:HD12	1.70	0.91
1:B:211:TYR:CE2	1:D:255:LEU:HD23	2.06	0.90
1:B:213:LEU:HD11	1:B:217:ARG:NE	1.86	0.90
1:A:130:ARG:NH1	1:A:265:THR:HG22	1.86	0.89
1:B:211:TYR:CD2	1:D:255:LEU:CD2	2.57	0.87
1:B:236:TYR:O	1:B:239:PRO:HD2	1.76	0.86
1:E:140:ARG:HH22	1:E:280:GLU:CD	1.83	0.86
1:E:140:ARG:NH2	1:E:280:GLU:OE2	2.08	0.85
1:C:201:ARG:HG3	1:E:83:ARG:HH11	1.43	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:VAL:HG21	1:B:260:GLU:CD	2.05	0.81
1:A:130:ARG:CZ	1:A:265:THR:HG22	2.11	0.80
1:B:211:TYR:HD2	1:D:255:LEU:CD2	1.92	0.80
1:B:211:TYR:HD2	1:D:255:LEU:HD23	1.46	0.79
1:A:130:ARG:HD2	1:A:265:THR:CG2	2.13	0.78
1:B:213:LEU:HD11	1:B:217:ARG:CZ	2.14	0.78
1:A:237:MET:HG2	1:E:50:ILE:HB	1.65	0.77
1:D:83:ARG:HG2	1:D:270:LEU:HD21	1.64	0.77
1:D:258:GLU:OE1	1:D:268:ASN:HB3	1.85	0.76
1:C:194:HIS:HA	1:E:91:ILE:HD13	1.68	0.75
1:E:135:THR:HG23	1:E:151:LEU:HD11	1.68	0.75
1:B:213:LEU:CD1	1:B:217:ARG:NE	2.50	0.74
1:E:130:ARG:NH1	1:E:269:ASP:OD2	2.19	0.74
1:A:149:GLU:OE2	1:B:291:PRO:HB3	1.88	0.74
1:C:91:ILE:HD11	1:D:197:GLY:HA3	1.70	0.73
1:C:285:ASP:O	1:D:159:ARG:NH2	2.22	0.72
1:A:37:ILE:O	1:A:41:SER:OG	2.08	0.71
1:D:268:ASN:O	1:D:268:ASN:ND2	2.21	0.71
1:A:234:LEU:HD23	1:E:51:HIS:O	1.90	0.71
1:A:258:GLU:HG2	1:A:268:ASN:HD22	1.56	0.71
1:E:208:PRO:HB3	1:E:212:THR:HG22	1.71	0.71
1:C:270:LEU:HD22	1:C:275:MET:HE1	1.72	0.71
1:B:213:LEU:C	1:B:213:LEU:HD12	2.17	0.70
1:B:213:LEU:HD12	1:B:213:LEU:O	1.91	0.70
1:B:104:LEU:HD11	1:B:177:LEU:HG	1.75	0.69
1:A:271:PRO:O	1:A:275:MET:HG3	1.93	0.69
1:C:39:ILE:HD11	1:C:239:PRO:HD3	1.74	0.69
1:A:214:ILE:HD13	1:B:251:SER:HB2	1.75	0.68
1:B:44:TRP:O	1:B:48:LEU:HD13	1.91	0.68
1:A:71:ARG:NH2	1:A:253:ASP:OD1	2.25	0.68
1:A:91:ILE:O	1:A:95:THR:HG23	1.93	0.68
1:D:71:ARG:NH2	1:D:253:ASP:OD1	2.28	0.67
1:A:52:LEU:HB3	1:A:229:ALA:HA	1.76	0.67
1:C:201:ARG:O	1:C:205:THR:HG22	1.94	0.67
1:A:130:ARG:HD2	1:A:265:THR:HG22	1.77	0.66
1:A:108:HIS:HB3	3:A:304:CL:CL	2.33	0.66
1:B:127:HIS:CD2	1:B:134:PRO:HA	2.31	0.66
1:D:143:PRO:HG2	1:D:146:ARG:HG2	1.77	0.66
1:E:43:GLN:O	1:E:46:GLU:HB2	1.95	0.65
1:A:258:GLU:CG	1:A:268:ASN:HD22	2.08	0.65
1:A:159:ARG:HH12	1:B:286:MET:HE3	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:LEU:HD22	1:A:275:MET:HE1	1.79	0.64
1:B:213:LEU:HD12	1:B:217:ARG:HD2	1.77	0.64
1:C:32:VAL:O	1:C:36:ILE:HG13	1.97	0.64
1:C:282:ASN:O	1:C:286:MET:HG3	1.97	0.64
1:A:258:GLU:HG2	1:A:268:ASN:ND2	2.12	0.64
1:C:25:ILE:HA	1:C:28:LEU:HD12	1.79	0.64
1:A:236:TYR:HB2	1:E:50:ILE:HG22	1.80	0.64
1:A:143:PRO:HG2	1:A:146:ARG:HG2	1.80	0.63
1:B:29:LEU:HA	1:B:32:VAL:HG23	1.79	0.63
1:A:93:GLU:HG2	1:A:118:LEU:HD22	1.81	0.63
1:B:207:VAL:HG21	1:B:260:GLU:OE2	1.97	0.63
1:A:52:LEU:HD12	1:A:228:PHE:HB3	1.80	0.62
1:B:60:LEU:HD13	1:D:245:ILE:HG12	1.80	0.62
1:B:90:LEU:HG	1:B:94:ARG:HD2	1.80	0.62
1:A:258:GLU:HA	1:A:258:GLU:OE1	2.00	0.61
1:C:277:ASN:HB3	1:C:281:ARG:HH22	1.65	0.61
1:C:201:ARG:HG3	1:E:83:ARG:NH1	2.15	0.61
1:C:212:THR:HG23	1:C:216:GLN:HG3	1.83	0.60
1:A:271:PRO:HG2	1:A:275:MET:HE3	1.84	0.60
1:E:213:LEU:HD11	1:E:217:ARG:CZ	2.30	0.60
1:A:130:ARG:CD	1:A:265:THR:CG2	2.80	0.60
1:A:149:GLU:HG2	1:B:291:PRO:HG3	1.83	0.60
1:A:271:PRO:HB2	1:A:274:ALA:HB3	1.84	0.59
1:B:52:LEU:HD23	1:D:234:LEU:HD22	1.84	0.59
1:D:71:ARG:HH12	1:D:212:THR:HG23	1.66	0.59
1:E:71:ARG:NH1	1:E:260:GLU:OE2	2.33	0.59
1:D:168:ILE:HG22	1:D:182:TYR:CD1	2.38	0.59
1:B:95:THR:HG22	1:B:98:ARG:HE	1.68	0.59
1:D:143:PRO:HB2	1:D:145:GLU:OE1	2.03	0.59
1:E:130:ARG:HH22	1:E:269:ASP:HA	1.68	0.59
1:E:140:ARG:NH2	1:E:280:GLU:CD	2.56	0.59
1:A:100:LEU:HD11	1:A:114:ILE:HD13	1.84	0.59
1:D:23:LYS:HA	1:D:26:PHE:HD2	1.66	0.58
1:E:126:LYS:HE3	1:E:273:ASN:OD1	2.03	0.58
1:C:135:THR:HG23	1:C:151:LEU:HD11	1.86	0.58
1:D:207:VAL:O	1:D:208:PRO:C	2.46	0.58
1:B:208:PRO:CB	1:D:268:ASN:HD21	2.17	0.58
1:C:255:LEU:HD11	1:D:210:ALA:HB3	1.86	0.58
1:E:271:PRO:O	1:E:275:MET:HB2	2.04	0.57
1:B:201:ARG:O	1:B:205:THR:HB	2.04	0.57
1:B:258:GLU:OE1	1:B:266:ALA:CB	2.53	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:281:ARG:HG2	1:D:155:MET:HE1	1.86	0.57
1:D:93:GLU:HB3	1:D:283:LEU:HD11	1.87	0.57
1:E:143:PRO:HG2	1:E:146:ARG:HG2	1.86	0.57
1:A:194:HIS:HA	1:B:91:ILE:HD13	1.87	0.57
1:B:101:ARG:HD3	1:B:290:HIS:HB3	1.86	0.57
1:A:159:ARG:NH1	1:B:286:MET:HE3	2.20	0.56
1:A:42:TYR:HA	1:A:45:TYR:HD2	1.70	0.56
1:B:139:ARG:HG2	1:B:147:VAL:HG11	1.86	0.56
1:C:150:ILE:HA	1:C:159:ARG:HG2	1.87	0.56
1:D:127:HIS:ND1	1:D:132:THR:OG1	2.37	0.56
1:B:93:GLU:HB3	1:B:283:LEU:HD11	1.88	0.56
1:B:27:ARG:HA	1:B:30:LEU:HB3	1.87	0.56
1:D:108:HIS:O	1:D:112:ARG:HG3	2.04	0.56
1:A:39:ILE:HD13	1:A:236:TYR:HD1	1.70	0.56
1:A:130:ARG:CD	1:A:265:THR:HG22	2.36	0.56
1:B:201:ARG:HG3	1:D:83:ARG:HE	1.71	0.56
1:C:255:LEU:HD11	1:D:210:ALA:CB	2.36	0.56
1:A:60:LEU:HD23	1:A:222:PHE:HD2	1.71	0.55
1:B:271:PRO:HG2	1:B:275:MET:HE3	1.88	0.55
1:E:71:ARG:NH1	1:E:212:THR:OG1	2.39	0.55
1:A:111:HIS:O	1:A:115:VAL:HG23	2.07	0.55
1:B:28:LEU:HD11	1:B:247:TYR:HB2	1.88	0.55
1:B:205:THR:HG21	1:D:83:ARG:HD3	1.88	0.55
1:E:139:ARG:HG2	1:E:144:GLU:OE1	2.06	0.55
1:E:139:ARG:HG3	1:E:147:VAL:HG11	1.89	0.55
1:B:115:VAL:O	1:B:119:VAL:HG23	2.08	0.54
1:D:149:GLU:OE2	1:D:159:ARG:HD2	2.07	0.54
1:A:22:SER:C	1:A:24:ILE:H	2.15	0.54
1:A:62:ILE:HG21	1:B:62:ILE:HG12	1.88	0.54
1:B:211:TYR:O	1:B:215:LEU:HB2	2.07	0.54
1:C:252:TRP:HZ3	1:D:215:LEU:HG	1.73	0.54
1:B:155:MET:HE3	1:D:94:ARG:NH2	2.23	0.54
1:B:208:PRO:HB3	1:D:268:ASN:HD21	1.72	0.53
1:B:60:LEU:HD23	1:B:222:PHE:HD2	1.73	0.53
1:D:71:ARG:HD2	1:D:71:ARG:C	2.34	0.53
1:A:210:ALA:HB1	1:B:255:LEU:HD11	1.91	0.53
1:C:115:VAL:HG21	1:C:287:THR:HG21	1.90	0.53
1:C:104:LEU:HD11	1:C:177:LEU:HG	1.90	0.53
1:C:240:PHE:O	1:C:243:VAL:HG12	2.09	0.53
1:C:40:ILE:HG13	1:C:41:SER:N	2.24	0.52
1:D:60:LEU:HD23	1:D:222:PHE:HD1	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ARG:NE	1:D:205:THR:OG1	2.39	0.52
1:E:140:ARG:NH2	1:E:280:GLU:OE1	2.42	0.52
1:E:208:PRO:HB3	1:E:212:THR:CG2	2.40	0.52
1:A:233:ASP:HB3	1:E:53:THR:OG1	2.09	0.52
1:B:211:TYR:CD2	1:D:255:LEU:HD21	2.43	0.52
1:E:207:VAL:N	1:E:208:PRO:HD2	2.25	0.52
1:C:237:MET:HG2	1:D:50:ILE:HA	1.92	0.52
1:C:275:MET:O	1:C:279:ILE:HG13	2.10	0.52
1:B:121:PHE:CE2	1:B:161:LEU:HG	2.46	0.51
1:C:35:SER:O	1:C:39:ILE:HG13	2.10	0.51
1:C:201:ARG:CG	1:E:83:ARG:HH11	2.20	0.51
1:B:113:ARG:O	1:B:116:SER:OG	2.26	0.51
1:D:72:ASN:HA	1:D:256:ALA:HB2	1.93	0.51
1:D:231:VAL:HA	1:D:238:THR:OG1	2.11	0.51
1:A:258:GLU:CG	1:A:268:ASN:ND2	2.71	0.50
1:C:146:ARG:HH11	1:C:146:ARG:HA	1.75	0.50
1:B:127:HIS:ND1	1:B:132:THR:OG1	2.42	0.50
1:C:238:THR:N	1:C:239:PRO:HD2	2.26	0.50
1:C:91:ILE:O	1:C:95:THR:HG23	2.11	0.50
1:E:86:TRP:CE2	1:E:272:LEU:HD11	2.46	0.50
1:B:213:LEU:HD11	1:B:217:ARG:NH1	2.25	0.50
1:C:83:ARG:HB2	1:C:270:LEU:HD21	1.94	0.50
1:B:258:GLU:OE1	1:B:266:ALA:HB1	2.11	0.49
1:E:100:LEU:HD11	1:E:114:ILE:HD13	1.94	0.49
1:A:27:ARG:HA	1:A:30:LEU:HB3	1.95	0.49
1:A:83:ARG:HD3	1:A:270:LEU:CD2	2.42	0.49
1:A:213:LEU:HD11	1:A:217:ARG:CZ	2.41	0.49
1:D:208:PRO:O	1:D:210:ALA:N	2.34	0.49
1:E:115:VAL:HG13	1:E:283:LEU:HB3	1.94	0.49
1:A:130:ARG:HD2	1:A:265:THR:HG23	1.95	0.49
1:A:43:GLN:HB2	1:A:44:TRP:HD1	1.78	0.49
1:C:83:ARG:HD3	1:C:270:LEU:HD22	1.94	0.49
1:B:134:PRO:HB2	1:B:138:LEU:HD12	1.95	0.49
1:C:168:ILE:HG22	1:C:182:TYR:CD1	2.48	0.49
1:B:71:ARG:NH2	1:B:253:ASP:OD1	2.42	0.48
1:D:268:ASN:HD22	1:D:268:ASN:C	2.19	0.48
1:C:207:VAL:HG13	1:E:76:TYR:HE1	1.78	0.48
1:A:230:LEU:HD21	1:E:56:PRO:HB3	1.96	0.48
1:A:98:ARG:HB2	1:A:286:MET:HE1	1.95	0.48
1:B:42:TYR:O	1:B:45:TYR:HB2	2.13	0.48
1:B:207:VAL:O	1:B:208:PRO:C	2.57	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:LEU:CD1	1:B:217:ARG:CG	2.92	0.48
1:A:31:ASN:ND2	1:A:223:CYS:O	2.47	0.48
1:D:71:ARG:NH1	1:D:212:THR:HG23	2.29	0.48
1:D:201:ARG:O	1:D:205:THR:HB	2.13	0.48
1:D:208:PRO:C	1:D:210:ALA:H	2.18	0.47
1:A:130:ARG:CD	1:A:265:THR:HG23	2.44	0.47
1:C:38:ALA:HB2	1:C:228:PHE:CD1	2.49	0.47
1:A:130:ARG:NE	1:A:265:THR:HG22	2.29	0.47
1:C:60:LEU:HB2	1:E:245:ILE:HG12	1.95	0.47
1:C:81:GLU:OE2	1:C:201:ARG:NH2	2.47	0.47
1:C:201:ARG:HG3	1:E:83:ARG:HE	1.79	0.47
1:A:207:VAL:O	1:A:208:PRO:C	2.57	0.47
1:A:255:LEU:HD13	1:E:211:TYR:CD2	2.49	0.47
1:B:101:ARG:NH1	1:B:288:GLY:C	2.73	0.47
1:D:278:THR:HG22	1:D:282:ASN:ND2	2.30	0.47
1:A:97:VAL:HG21	1:A:283:LEU:HD22	1.96	0.47
1:B:104:LEU:HB3	1:B:107:GLU:HG3	1.97	0.47
1:E:71:ARG:HD2	1:E:71:ARG:C	2.40	0.47
1:B:215:LEU:HD21	1:D:252:TRP:CE3	2.50	0.46
1:B:168:ILE:HG22	1:B:182:TYR:CD1	2.50	0.46
1:C:236:TYR:C	1:C:238:THR:H	2.23	0.46
1:A:153:SER:OG	1:B:285:ASP:OD2	2.16	0.46
1:E:81:GLU:OE1	1:E:202:LEU:HD11	2.15	0.46
1:B:140:ARG:NH1	1:B:141:LEU:HD21	2.29	0.46
1:C:55:ALA:HB3	1:C:56:PRO:HD3	1.97	0.46
1:C:62:ILE:HG12	1:D:62:ILE:HG21	1.97	0.46
1:A:234:LEU:CD2	1:E:51:HIS:O	2.60	0.46
1:B:85:LEU:HB3	1:B:195:VAL:HG13	1.98	0.46
1:C:171:LEU:HD23	1:C:171:LEU:HA	1.81	0.46
1:A:45:TYR:HD1	1:A:50:ILE:O	1.99	0.46
1:D:127:HIS:NE2	1:D:137:ASP:OD2	2.33	0.46
1:A:194:HIS:CE1	1:B:191:GLU:OE2	2.69	0.46
1:C:157:THR:HB	1:C:196:LEU:HD22	1.98	0.46
1:C:37:ILE:O	1:C:41:SER:OG	2.16	0.45
1:E:130:ARG:HH12	1:E:269:ASP:CG	2.24	0.45
1:A:31:ASN:O	1:A:35:SER:N	2.49	0.45
1:B:24:ILE:HD13	1:B:27:ARG:HH22	1.81	0.45
1:D:204:THR:O	1:D:206:PRO:HD3	2.17	0.45
1:B:207:VAL:CG2	1:B:260:GLU:CD	2.84	0.45
1:C:207:VAL:HG13	1:E:76:TYR:CE1	2.52	0.45
1:D:33:LEU:HA	1:D:36:ILE:HG12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:MET:HE3	1:E:52:LEU:HD21	1.98	0.45
1:D:191:GLU:O	1:D:195:VAL:HG23	2.17	0.45
1:A:44:TRP:O	1:A:48:LEU:HD13	2.17	0.45
1:E:107:GLU:CD	1:E:176:LYS:HD3	2.42	0.45
1:C:60:LEU:HD21	1:E:244:PHE:CE2	2.51	0.45
1:E:208:PRO:HB2	1:E:210:ALA:N	2.32	0.45
1:C:221:LEU:HD23	1:E:244:PHE:CE1	2.52	0.45
1:D:270:LEU:HA	1:D:270:LEU:HD23	1.74	0.45
1:E:45:TYR:HB2	1:E:50:ILE:HG13	1.99	0.45
1:A:129:LEU:HD21	1:A:199:CYS:HB3	1.99	0.44
1:B:117:TYR:CZ	1:B:163:LEU:HB3	2.53	0.44
1:C:76:TYR:OH	1:D:207:VAL:HG22	2.16	0.44
1:C:89:VAL:CG2	1:C:195:VAL:HG11	2.47	0.44
1:D:109:ASP:HB3	1:D:113:ARG:HH12	1.82	0.44
1:A:44:TRP:CD1	1:A:44:TRP:N	2.85	0.44
1:B:89:VAL:HG23	1:B:195:VAL:HG11	1.98	0.44
1:B:119:VAL:HG11	1:B:280:GLU:HG3	2.00	0.44
1:B:215:LEU:HD21	1:D:252:TRP:CZ3	2.52	0.44
1:C:236:TYR:O	1:C:239:PRO:CD	2.54	0.44
1:D:22:SER:O	1:D:25:ILE:HG22	2.18	0.44
1:E:103:ILE:O	1:E:105:PRO:HD3	2.18	0.44
1:B:143:PRO:O	1:B:147:VAL:HG23	2.18	0.44
1:C:240:PHE:O	1:C:241:VAL:C	2.60	0.44
1:D:50:ILE:O	1:D:50:ILE:HG13	2.18	0.44
1:B:230:LEU:HB3	1:B:234:LEU:HD12	1.99	0.43
1:A:36:ILE:O	1:A:40:ILE:HG12	2.18	0.43
1:B:213:LEU:CD1	1:B:217:ARG:CZ	2.91	0.43
1:D:68:LEU:HB3	1:D:252:TRP:HD1	1.82	0.43
1:C:100:LEU:HD11	1:C:114:ILE:HD13	1.99	0.43
1:D:272:LEU:HA	1:D:272:LEU:HD23	1.80	0.43
1:A:93:GLU:HB3	1:A:283:LEU:HD11	2.00	0.43
1:B:26:PHE:HD1	1:B:27:ARG:HG3	1.83	0.43
1:B:191:GLU:O	1:B:195:VAL:HG23	2.17	0.43
1:C:181:THR:HA	1:C:184:LEU:HG	1.99	0.43
1:C:202:LEU:HD23	1:C:202:LEU:HA	1.66	0.43
1:E:38:ALA:O	1:E:42:TYR:HB2	2.19	0.43
1:C:83:ARG:HG2	1:D:205:THR:HG21	2.01	0.43
1:A:52:LEU:HB3	1:A:229:ALA:CA	2.47	0.43
1:A:110:ALA:HB1	1:A:171:LEU:HD11	2.00	0.43
1:A:130:ARG:HH11	1:A:265:THR:HG22	1.73	0.43
1:C:271:PRO:O	1:C:275:MET:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:168:ILE:HG22	1:D:182:TYR:HD1	1.79	0.43
1:B:155:MET:HE3	1:D:94:ARG:HH21	1.83	0.42
1:E:121:PHE:CE1	1:E:157:THR:HG22	2.54	0.42
1:C:236:TYR:C	1:C:238:THR:N	2.77	0.42
1:A:49:GLY:O	1:B:235:HIS:HB2	2.18	0.42
1:B:211:TYR:HE2	1:D:255:LEU:HD23	1.77	0.42
1:D:143:PRO:HG2	1:D:146:ARG:CG	2.47	0.42
1:A:52:LEU:HD12	1:A:228:PHE:CB	2.48	0.42
1:A:57:PHE:HD2	1:A:226:LEU:HD12	1.84	0.42
1:B:286:MET:O	1:B:289:GLN:HB3	2.20	0.42
1:D:145:GLU:OE1	1:D:145:GLU:N	2.44	0.42
1:A:211:TYR:HA	1:B:255:LEU:HD21	2.00	0.42
1:A:42:TYR:O	1:A:45:TYR:HB2	2.19	0.42
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.78	0.42
1:C:192:LEU:HD23	1:C:192:LEU:HA	1.77	0.42
1:B:101:ARG:HH11	1:B:288:GLY:C	2.28	0.42
1:C:42:TYR:HA	1:C:45:TYR:CD1	2.55	0.42
1:C:252:TRP:CH2	1:D:67:PHE:HB3	2.54	0.42
1:A:83:ARG:HD3	1:A:270:LEU:HD21	2.02	0.42
1:B:270:LEU:HD23	1:B:270:LEU:HA	1.86	0.41
1:C:79:PHE:HB2	1:C:259:ALA:HB1	2.02	0.41
1:C:290:HIS:HA	1:C:291:PRO:HD2	1.83	0.41
1:D:129:LEU:HD21	1:D:199:CYS:HB3	2.02	0.41
1:E:213:LEU:HD21	1:E:217:ARG:NH1	2.35	0.41
1:E:269:ASP:OD1	1:E:273:ASN:HB2	2.19	0.41
1:C:39:ILE:HD13	1:C:236:TYR:HA	2.01	0.41
1:E:71:ARG:NH2	1:E:253:ASP:OD1	2.51	0.41
1:E:208:PRO:O	1:E:211:TYR:HB2	2.20	0.41
1:A:60:LEU:HD13	1:B:245:ILE:HG12	2.02	0.41
1:C:240:PHE:N	1:C:240:PHE:CD1	2.86	0.41
1:D:208:PRO:C	1:D:210:ALA:N	2.78	0.41
1:E:90:LEU:HD12	1:E:279:ILE:HB	2.02	0.41
1:A:135:THR:HG23	1:A:151:LEU:HD11	2.01	0.41
1:A:262:PRO:O	1:A:270:LEU:HB2	2.21	0.41
1:B:101:ARG:HD3	1:B:290:HIS:CB	2.49	0.41
1:B:161:LEU:HD23	1:B:161:LEU:HA	1.89	0.41
1:B:208:PRO:O	1:B:211:TYR:HB2	2.21	0.41
1:C:284:LEU:HA	1:C:284:LEU:HD23	1.86	0.41
1:D:55:ALA:HB3	1:D:56:PRO:HD3	2.02	0.41
1:D:104:LEU:HD21	1:D:176:LYS:O	2.20	0.41
1:A:22:SER:C	1:A:24:ILE:N	2.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:LEU:HD12	1:B:30:LEU:O	2.20	0.41
1:C:139:ARG:HG3	1:C:147:VAL:HG11	2.03	0.41
1:A:168:ILE:HD12	1:A:168:ILE:HG23	1.80	0.41
1:B:52:LEU:HD23	1:B:52:LEU:HA	1.89	0.41
1:B:168:ILE:HG22	1:B:182:TYR:HD1	1.86	0.41
1:C:100:LEU:HD11	1:C:114:ILE:CD1	2.50	0.41
1:C:30:LEU:O	1:C:34:MET:HG2	2.21	0.41
1:C:207:VAL:O	1:C:208:PRO:C	2.64	0.41
1:A:45:TYR:CD1	1:A:50:ILE:HG13	2.55	0.41
1:B:89:VAL:CG2	1:B:195:VAL:HG11	2.51	0.41
1:B:171:LEU:HA	1:B:171:LEU:HD23	1.80	0.41
1:B:236:TYR:C	1:B:238:THR:H	2.29	0.41
1:C:103:ILE:HD12	1:C:185:MET:HE3	2.03	0.41
1:C:129:LEU:HD23	1:C:129:LEU:HA	1.82	0.41
1:C:168:ILE:HG23	1:C:168:ILE:HD12	1.73	0.41
1:E:108:HIS:HB3	3:E:305:CL:CL	2.57	0.41
1:E:150:ILE:HA	1:E:159:ARG:HG2	2.02	0.41
1:A:72:ASN:OD1	1:A:256:ALA:HB2	2.21	0.40
1:A:200:GLU:O	1:A:204:THR:HG23	2.20	0.40
1:C:36:ILE:HG13	1:C:36:ILE:H	1.67	0.40
1:E:33:LEU:O	1:E:36:ILE:HG22	2.21	0.40
1:A:85:LEU:HD22	1:A:195:VAL:HA	2.03	0.40
1:C:104:LEU:HD12	1:C:171:LEU:HD13	2.03	0.40
1:A:42:TYR:O	1:A:43:GLN:C	2.65	0.40
1:A:42:TYR:HD1	1:A:45:TYR:CD2	2.39	0.40
1:B:42:TYR:CD1	1:B:45:TYR:CD2	3.09	0.40
1:E:279:ILE:HD12	1:E:279:ILE:HA	1.87	0.40
1:B:129:LEU:HD11	1:B:199:CYS:HB3	2.03	0.40
1:B:140:ARG:O	1:B:140:ARG:HG3	2.19	0.40
1:E:177:LEU:HD23	1:E:177:LEU:HA	1.89	0.40
1:B:25:ILE:HA	1:B:28:LEU:HB2	2.03	0.40
1:D:215:LEU:HD23	1:D:215:LEU:HA	1.89	0.40
1:E:272:LEU:HD12	1:E:272:LEU:HA	1.79	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	266/297 (90%)	251 (94%)	15 (6%)	0	100	100
1	B	268/297 (90%)	248 (92%)	20 (8%)	0	100	100
1	C	267/297 (90%)	249 (93%)	18 (7%)	0	100	100
1	D	266/297 (90%)	248 (93%)	16 (6%)	2 (1%)	16	44
1	E	268/297 (90%)	257 (96%)	11 (4%)	0	100	100
All	All	1335/1485 (90%)	1253 (94%)	80 (6%)	2 (0%)	48	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	209	PHE
1	D	208	PRO

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	232/259 (90%)	232 (100%)	0	100	100
1	B	234/259 (90%)	232 (99%)	2 (1%)	70	76
1	C	232/259 (90%)	231 (100%)	1 (0%)	84	83
1	D	232/259 (90%)	230 (99%)	2 (1%)	70	76
1	E	234/259 (90%)	234 (100%)	0	100	100
All	All	1164/1295 (90%)	1159 (100%)	5 (0%)	84	83

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	213	LEU
1	B	258	GLU
1	C	265	THR
1	D	257	GLU
1	D	258	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	268	ASN
1	A	289	GLN
1	B	273	ASN
1	B	290	HIS
1	C	102	ASN
1	C	166	ASN
1	C	273	ASN
1	D	31	ASN
1	D	47	GLN
1	D	84	ASN
1	D	216	GLN
1	D	268	ASN
1	D	289	GLN
1	E	72	ASN
1	E	99	GLN
1	E	166	ASN
1	E	282	ASN

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 [i](#)

Of 23 ligands modelled in this entry, 22 are monoatomic - leaving 1 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	$\# Z > 2$
4	ACY	C	304	-	3,3,3	2.81	1 (33%)	3,3,3	1.46	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	304	ACY	CH3-C	4.27	1.65	1.49

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

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	268/297 (90%)	0.52	24 (8%) 15 14	58, 90, 150, 173	0
1	B	270/297 (90%)	0.68	41 (15%) 5 7	58, 100, 161, 213	0
1	C	269/297 (90%)	0.67	36 (13%) 7 8	60, 93, 158, 190	0
1	D	268/297 (90%)	1.39	70 (26%) 1 3	64, 103, 151, 204	0
1	E	270/297 (90%)	0.74	44 (16%) 4 6	56, 89, 168, 228	0
All	All	1345/1485 (90%)	0.80	215 (15%) 5 6	56, 96, 157, 228	0

All (215) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	GLU	16.6
1	D	145	GLU	15.2
1	D	265	THR	14.9
1	D	233	ASP	12.8
1	C	109	ASP	11.3
1	D	131	LYS	10.2
1	D	146	ARG	9.6
1	C	106	ALA	9.5
1	D	166	ASN	9.5
1	D	173	GLU	9.1
1	D	236	TYR	8.8
1	B	47	GLN	8.6
1	A	175	GLY	8.4
1	E	47	GLN	8.4
1	B	51	HIS	8.1
1	B	49	GLY	8.0
1	B	220	TYR	7.4
1	D	235	HIS	7.3
1	C	47	GLN	7.2
1	D	170	GLN	7.0

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Mol	Chain	Res	Type	RSRZ
1	D	42	TYR	6.3
1	D	46	GLU	6.1
1	D	43	GLN	6.0
1	D	47	GLN	6.0
1	D	232	GLY	5.9
1	D	130	ARG	5.8
1	D	106	ALA	5.6
1	C	24	ILE	5.5
1	D	149	GLU	5.5
1	C	107	GLU	5.3
1	E	170	GLN	5.3
1	B	53	THR	5.3
1	C	208	PRO	5.3
1	B	288	GLY	5.2
1	C	184	LEU	5.2
1	A	174	ALA	5.2
1	E	167	GLU	5.1
1	B	216	GLN	5.1
1	E	235	HIS	5.1
1	D	211	TYR	5.1
1	E	265	THR	5.0
1	A	176	LYS	5.0
1	D	116	SER	4.9
1	E	173	GLU	4.9
1	D	167	GLU	4.8
1	D	209	PHE	4.8
1	A	140	ARG	4.7
1	C	105	PRO	4.7
1	B	26	PHE	4.7
1	B	50	ILE	4.6
1	E	174	ALA	4.6
1	D	153	SER	4.6
1	C	48	LEU	4.5
1	E	266	ALA	4.5
1	D	129	LEU	4.5
1	D	163	LEU	4.4
1	B	48	LEU	4.4
1	E	168	ILE	4.4
1	A	173	GLU	4.3
1	B	207	VAL	4.2
1	D	266	ALA	4.2
1	E	36	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	44	TRP	4.1
1	E	209	PHE	4.1
1	D	143	PRO	4.1
1	B	73	SER	4.0
1	D	174	ALA	4.0
1	B	287	THR	3.8
1	E	42	TYR	3.8
1	C	213	LEU	3.7
1	E	45	TYR	3.7
1	B	40	ILE	3.7
1	B	208	PRO	3.6
1	D	208	PRO	3.6
1	E	25	ILE	3.6
1	E	39	ILE	3.6
1	C	25	ILE	3.6
1	E	48	LEU	3.5
1	D	107	GLU	3.5
1	D	184	LEU	3.5
1	D	117	TYR	3.5
1	A	105	PRO	3.5
1	C	26	PHE	3.4
1	D	187	ASN	3.4
1	D	113	ARG	3.4
1	A	103	ILE	3.4
1	B	215	LEU	3.4
1	A	209	PHE	3.3
1	C	260	GLU	3.3
1	D	207	VAL	3.3
1	E	216	GLN	3.3
1	B	52	LEU	3.3
1	E	254	SER	3.3
1	C	207	VAL	3.3
1	D	243	VAL	3.3
1	B	211	TYR	3.2
1	B	212	THR	3.2
1	D	224	THR	3.2
1	D	264	GLY	3.2
1	A	32	VAL	3.2
1	B	228	PHE	3.2
1	C	145	GLU	3.1
1	D	139	ARG	3.1
1	E	110	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	255	LEU	3.1
1	D	234	LEU	3.1
1	E	251	SER	3.1
1	B	39	ILE	3.1
1	A	106	ALA	3.1
1	B	289	GLN	3.0
1	E	176	LYS	3.0
1	C	211	TYR	3.0
1	B	24	ILE	3.0
1	C	259	ALA	3.0
1	D	31	ASN	3.0
1	D	27	ARG	2.9
1	D	28	LEU	2.9
1	D	180	ILE	2.9
1	E	222	PHE	2.9
1	B	224	THR	2.8
1	D	246	SER	2.8
1	D	206	PRO	2.8
1	D	186	ASP	2.8
1	A	204	THR	2.8
1	E	268	ASN	2.8
1	E	46	GLU	2.8
1	A	35	SER	2.7
1	D	273	ASN	2.7
1	D	110	ALA	2.7
1	C	209	PHE	2.7
1	C	254	SER	2.7
1	B	233	ASP	2.7
1	A	102	ASN	2.7
1	E	232	GLY	2.6
1	D	223	CYS	2.6
1	E	231	VAL	2.6
1	E	253	ASP	2.6
1	C	187	ASN	2.6
1	D	176	LYS	2.6
1	C	46	GLU	2.6
1	A	224	THR	2.6
1	A	104	LEU	2.6
1	C	95	THR	2.6
1	B	93	GLU	2.6
1	E	113	ARG	2.6
1	E	172	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	214	ILE	2.6
1	C	227	PRO	2.6
1	D	39	ILE	2.5
1	C	252	TRP	2.5
1	E	37	ILE	2.5
1	A	177	LEU	2.5
1	B	225	LEU	2.5
1	B	35	SER	2.5
1	C	212	THR	2.5
1	D	154	SER	2.5
1	C	291	PRO	2.5
1	D	32	VAL	2.5
1	B	253	ASP	2.4
1	C	180	ILE	2.4
1	D	30	LEU	2.4
1	D	250	LEU	2.4
1	E	219	VAL	2.4
1	C	110	ALA	2.4
1	B	209	PHE	2.4
1	C	51	HIS	2.4
1	D	183	GLY	2.4
1	D	210	ALA	2.4
1	A	223	CYS	2.4
1	E	40	ILE	2.4
1	D	179	ASP	2.3
1	C	29	LEU	2.3
1	C	99	GLN	2.3
1	A	36	ILE	2.3
1	B	34	MET	2.3
1	E	205	THR	2.3
1	D	228	PHE	2.3
1	E	220	TYR	2.3
1	E	24	ILE	2.2
1	B	286	MET	2.2
1	D	268	ASN	2.2
1	E	64	ILE	2.2
1	B	77	SER	2.2
1	D	225	LEU	2.2
1	E	33	LEU	2.2
1	A	205	THR	2.2
1	A	30	LEU	2.2
1	A	48	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	249	PHE	2.2
1	B	257	GLU	2.1
1	D	144	GLU	2.1
1	A	44	TRP	2.1
1	C	44	TRP	2.1
1	E	53	THR	2.1
1	B	27	ARG	2.1
1	D	112	ARG	2.1
1	D	227	PRO	2.1
1	A	70	PHE	2.1
1	C	268	ASN	2.1
1	C	204	THR	2.1
1	A	107	GLU	2.1
1	B	260	GLU	2.1
1	B	122	SER	2.1
1	E	228	PHE	2.1
1	D	48	LEU	2.1
1	C	181	THR	2.1
1	D	220	TYR	2.1
1	E	210	ALA	2.1
1	E	264	GLY	2.1
1	E	206	PRO	2.1
1	B	285	ASP	2.0
1	D	135	THR	2.0
1	D	182	TYR	2.0
1	E	204	THR	2.0
1	D	193	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ACY	C	304	4/4	0.72	0.32	67,85,87,90	0
2	ZN	B	301	1/1	0.84	0.14	187,187,187,187	0
2	ZN	D	302	1/1	0.92	0.07	123,123,123,123	0
3	CL	D	304	1/1	0.94	0.10	119,119,119,119	0
2	ZN	B	304	1/1	0.95	0.12	155,155,155,155	0
2	ZN	A	303	1/1	0.95	0.07	126,126,126,126	0
2	ZN	D	303	1/1	0.96	0.07	156,156,156,156	0
3	CL	D	305	1/1	0.97	0.08	121,121,121,121	0
3	CL	E	304	1/1	0.97	0.07	81,81,81,81	0
2	ZN	A	301	1/1	0.97	0.09	89,89,89,89	0
3	CL	C	305	1/1	0.98	0.17	83,83,83,83	0
2	ZN	B	302	1/1	0.98	0.15	85,85,85,85	0
2	ZN	C	302	1/1	0.98	0.09	91,91,91,91	0
2	ZN	E	303	1/1	0.98	0.11	96,96,96,96	0
3	CL	E	305	1/1	0.98	0.07	91,91,91,91	0
3	CL	A	304	1/1	0.98	0.07	107,107,107,107	0
2	ZN	E	301	1/1	0.99	0.08	83,83,83,83	0
2	ZN	E	302	1/1	0.99	0.02	88,88,88,88	0
2	ZN	B	303	1/1	0.99	0.08	84,84,84,84	0
2	ZN	A	302	1/1	0.99	0.09	101,101,101,101	0
2	ZN	C	301	1/1	0.99	0.10	90,90,90,90	0
2	ZN	C	303	1/1	1.00	0.03	104,104,104,104	0
2	ZN	D	301	1/1	1.00	0.02	105,105,105,105	0

6.5 Other polymers ⓘ

There are no such residues in this entry.