



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:35 AM GMT

PDB ID : 2AN6
Title : Protein-peptide complex
Authors : House, C.M.; Hancock, N.C.; Moller, A.; Cromer, B.A.; Fedorov, V.; Bowtell, D.D.L.; Parker, M.W.; Polekhina, G.
Deposited on : 2005-08-11
Resolution : 3.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

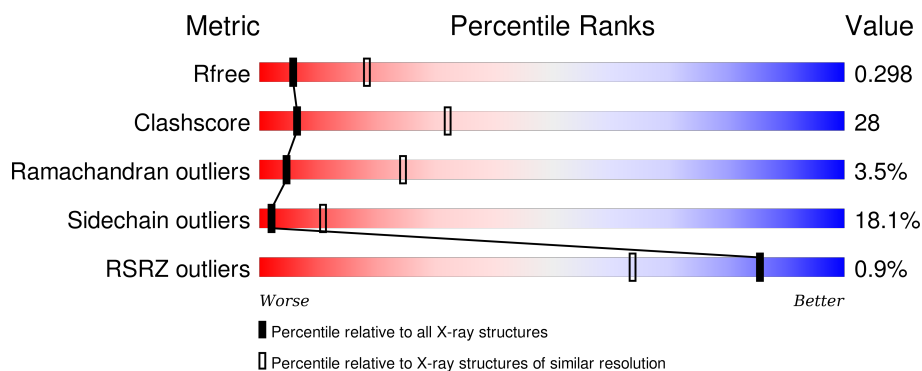
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	191	<div> <div>47%</div> <div>42%</div> <div>12%</div> </div>
1	B	191	<div> <div>%</div> <div>51%</div> <div>38%</div> <div>10%</div> <div>.</div> </div>
1	C	191	<div> <div>2%</div> <div>56%</div> <div>36%</div> <div>6%</div> <div>.</div> </div>
1	D	191	<div> <div>%</div> <div>47%</div> <div>39%</div> <div>13%</div> <div>.</div> </div>
2	E	24	<div> <div>25%</div> <div>21%</div> <div>54%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	24	 25% 21% 54%
2	G	24	 33% 8% 4% 54%
2	H	24	 4% 17% 13% 17% 54%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6372 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 Ubiquitin ligase SIAH1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1506	953	258	279	16			
1	B	191	Total	C	N	O	S	0	0	0
			1506	953	258	279	16			
1	C	191	Total	C	N	O	S	0	0	0
			1506	953	258	279	16			
1	D	191	Total	C	N	O	S	0	0	0
			1506	953	258	279	16			

- Molecule 2 is a protein called peptide from Phyllopod.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	11	Total	C	N	O	S	0	0	0
			85	55	17	12	1			
2	F	11	Total	C	N	O	S	0	0	0
			85	55	17	12	1			
2	G	11	Total	C	N	O	S	0	0	0
			85	55	17	12	1			
2	H	11	Total	C	N	O	S	0	0	0
			85	55	17	12	1			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

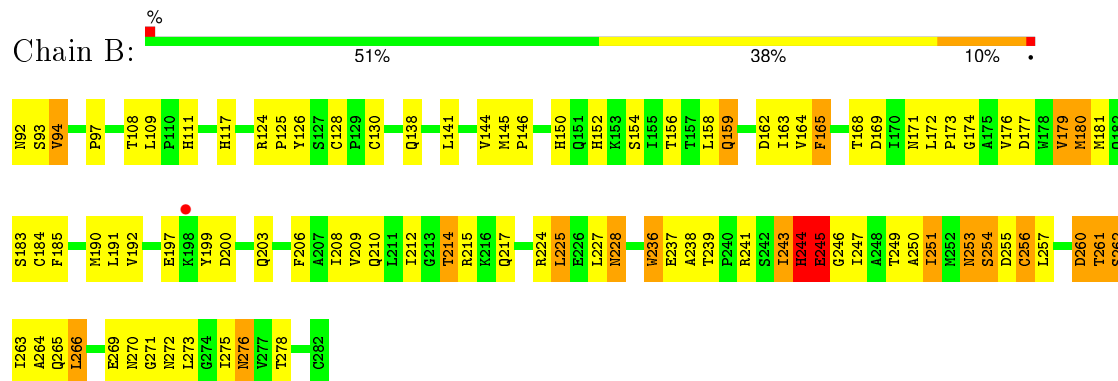
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

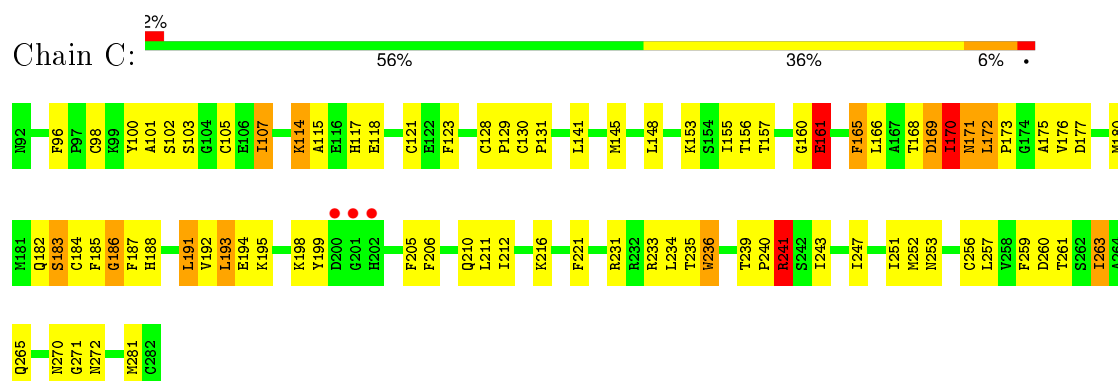
• Molecule 1: Ubiquitin ligase SIAH1A



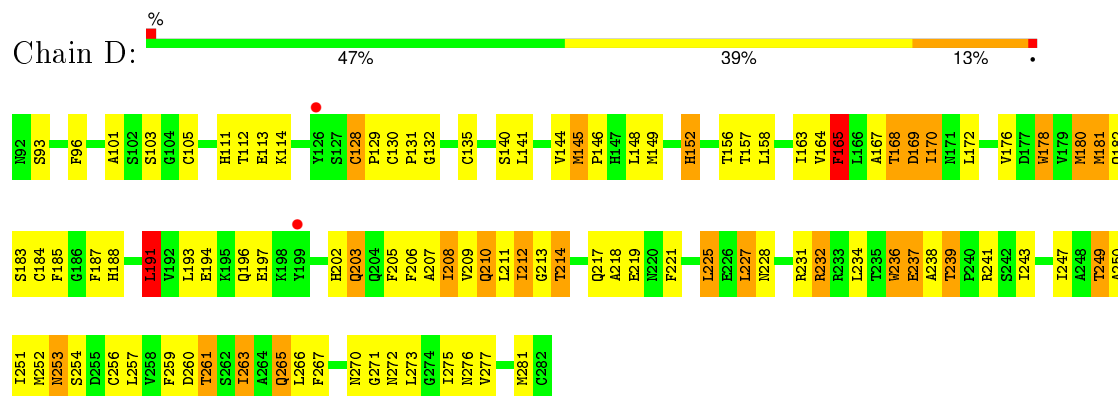
• Molecule 1: Ubiquitin ligase SIAH1A



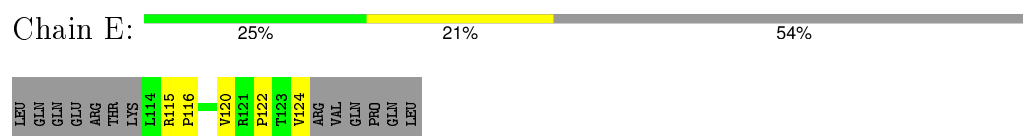
• Molecule 1: Ubiquitin ligase SIAH1A



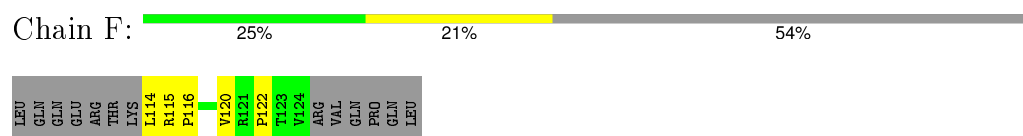
- Molecule 1: Ubiquitin ligase SIAH1A



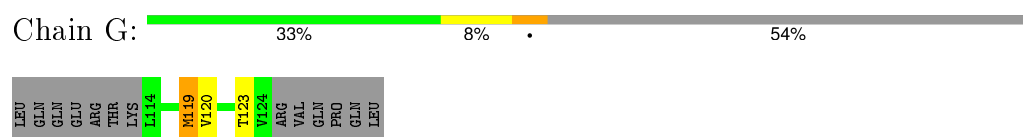
- Molecule 2: peptide from Phyllopod



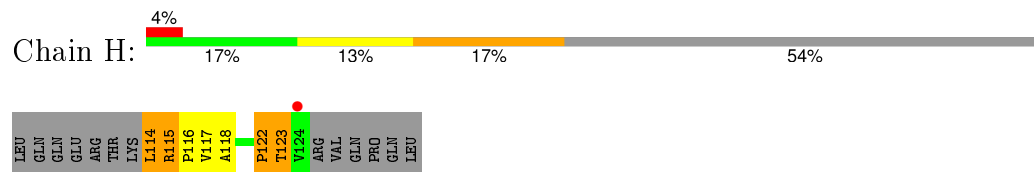
- Molecule 2: peptide from Phyllopod



- Molecule 2: peptide from Phyllopod



- Molecule 2: peptide from Phyllopod



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.30Å 100.00Å 103.40Å 90.00° 104.30° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 100.20 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-3.00) 98.5 (100.20-2.99)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.237 , 0.305 0.233 , 0.298	Depositor DCC
R_{free} test set	1257 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	73.8	Xtriage
Anisotropy	0.638	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 82.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 25343 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6372	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.73% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.81	0/1543	0.88	1/2088 (0.0%)
1	B	0.79	1/1543 (0.1%)	0.86	2/2088 (0.1%)
1	C	0.83	0/1543	0.86	0/2088
1	D	0.79	2/1543 (0.1%)	0.86	5/2088 (0.2%)
2	E	0.82	0/86	0.99	0/117
2	F	0.71	0/86	0.83	0/117
2	G	0.82	0/86	0.84	0/117
2	H	0.94	0/86	0.96	0/117
All	All	0.81	3/6516 (0.0%)	0.87	8/8820 (0.1%)

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
1	D	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	236	TRP	CE3-CZ3	7.52	1.51	1.38
1	D	165	PHE	CE2-CZ	5.13	1.47	1.37
1	B	236	TRP	CA-CB	5.11	1.65	1.53

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	225	LEU	CB-CG-CD1	-6.52	99.92	111.00
1	D	148	LEU	CA-CB-CG	6.05	129.22	115.30
1	B	225	LEU	CB-CG-CD1	-5.81	101.13	111.00
1	D	191	LEU	CA-CB-CG	5.66	128.31	115.30
1	D	225	LEU	CA-CB-CG	5.61	128.19	115.30
1	B	164	VAL	CB-CA-C	-5.52	100.91	111.40
1	D	165	PHE	N-CA-C	-5.44	96.32	111.00
1	A	211	LEU	CB-CG-CD2	-5.00	102.50	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	102	SER	Peptide
1	D	191	LEU	Peptide

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	1506	0	1451	87	0
1	B	1506	0	1454	91	0
1	C	1506	0	1451	75	0
1	D	1506	0	1454	108	0
2	E	85	0	98	5	0
2	F	85	0	98	7	0
2	G	85	0	98	3	0
2	H	85	0	98	6	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	1	0
All	All	6372	0	6202	346	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 28.

All (346) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:212:ILE:HA	1:C:243:ILE:HD11	1.36	1.08
1:B:168:THR:HG22	2:F:120:VAL:O	1.56	1.06
1:C:107:ILE:H	1:C:107:ILE:HD13	1.20	1.04
1:B:162:ASP:OD1	1:B:278:THR:HG22	1.58	1.00
1:A:254:SER:O	1:B:263:ILE:HD12	1.68	0.94
1:A:214:THR:OG1	1:A:217:GLN:HG3	1.69	0.92
1:B:253:ASN:O	1:B:254:SER:HB2	1.72	0.87
1:B:210:GLN:NE2	1:B:247:ILE:HG12	1.90	0.85
1:D:165:PHE:HE2	1:D:167:ALA:HB2	1.42	0.84
1:C:107:ILE:N	1:C:107:ILE:HD13	1.93	0.82
1:B:243:ILE:HD13	1:B:243:ILE:H	1.43	0.81
1:D:214:THR:H	1:D:217:GLN:HE21	1.29	0.79
1:D:145:MET:HE3	1:D:145:MET:HA	1.65	0.78
1:A:259:PHE:HB2	1:A:263:ILE:HG21	1.66	0.77
1:A:195:LYS:HD2	1:A:205:PHE:CE2	2.20	0.77
1:D:165:PHE:CE2	1:D:167:ALA:HB2	2.19	0.77
1:B:228:ASN:HD22	1:B:228:ASN:N	1.82	0.76
1:B:236:TRP:HB3	1:B:257:LEU:HB2	1.68	0.75
1:B:179:VAL:HG13	1:B:192:VAL:HG22	1.67	0.75
1:D:206:PHE:HB3	1:D:251:ILE:HD12	1.67	0.75
1:A:129:PRO:HG2	1:A:148:LEU:HD21	1.70	0.74
1:B:162:ASP:CG	1:B:278:THR:HG22	2.09	0.72
1:B:236:TRP:CZ2	1:B:238:ALA:HB2	2.24	0.72
1:D:163:ILE:HG22	1:D:164:VAL:N	2.05	0.72
1:D:206:PHE:HB3	1:D:251:ILE:CD1	2.20	0.72
1:C:193:LEU:HD12	1:C:194:GLU:N	2.05	0.72
1:C:183:SER:HB2	1:C:188:HIS:ND1	2.04	0.71
1:D:158:LEU:HD23	1:D:163:ILE:CG2	2.21	0.71
1:C:261:THR:O	1:C:265:GLN:HG2	1.90	0.71
1:D:253:ASN:ND2	1:D:253:ASN:H	1.87	0.70
1:D:253:ASN:H	1:D:253:ASN:HD22	1.38	0.70
1:D:232:ARG:HH11	1:D:232:ARG:HG3	1.55	0.70
1:A:176:VAL:HG12	1:A:177:ASP:N	2.06	0.70
1:D:145:MET:CE	1:D:145:MET:HA	2.20	0.70
1:B:247:ILE:HD13	1:B:251:ILE:HD11	1.73	0.69
1:B:158:LEU:HD21	2:F:116:PRO:HB2	1.72	0.69
1:D:197:GLU:HB2	1:D:203:GLN:HG3	1.74	0.69
1:D:168:THR:HA	1:D:272:ASN:OD1	1.93	0.69
1:D:209:VAL:HG23	1:D:225:LEU:HD21	1.76	0.68
1:A:233:ARG:HG2	1:B:237:GLU:OE1	1.93	0.67
1:A:232:ARG:HB2	1:B:236:TRP:CH2	2.30	0.67
1:A:101:ALA:HA	1:A:105:CYS:HB3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:VAL:HG12	1:A:177:ASP:H	1.60	0.67
1:B:276:ASN:HD22	1:B:276:ASN:N	1.92	0.66
1:A:204:GLN:OE1	1:A:206:PHE:CZ	2.49	0.66
1:C:103:SER:O	1:C:123:PHE:CD2	2.49	0.66
1:C:168:THR:HB	2:G:120:VAL:O	1.96	0.66
1:D:135:CYS:SG	1:D:152:HIS:HE1	2.17	0.66
1:A:162:ASP:CG	1:A:278:THR:HG22	2.15	0.66
1:A:204:GLN:OE1	1:A:206:PHE:HZ	1.79	0.65
1:C:212:ILE:CA	1:C:243:ILE:HD11	2.22	0.65
1:D:260:ASP:O	1:D:263:ILE:HG13	1.97	0.65
1:D:165:PHE:HD2	1:D:165:PHE:C	2.00	0.64
1:B:243:ILE:CD1	1:B:243:ILE:H	2.08	0.64
1:D:227:LEU:HD21	1:D:257:LEU:HD21	1.78	0.64
1:A:99:LYS:HB2	1:A:118:GLU:OE2	1.97	0.64
1:D:170:ILE:HD12	1:D:273:LEU:HB2	1.78	0.64
1:B:159:GLN:NE2	1:C:96:PHE:HE2	1.96	0.64
1:B:263:ILE:C	1:B:265:GLN:H	2.01	0.64
1:B:241:ARG:NH1	1:B:245:GLU:OE2	2.31	0.63
1:C:107:ILE:H	1:C:107:ILE:CD1	1.93	0.63
1:A:114:LYS:O	1:A:118:GLU:HB2	1.99	0.63
1:D:180:MET:HG3	1:D:181:MET:N	2.14	0.62
1:D:232:ARG:HH11	1:D:232:ARG:CG	2.12	0.62
1:D:135:CYS:SG	1:D:152:HIS:CE1	2.93	0.62
1:C:185:PHE:O	1:C:187:PHE:HD1	1.83	0.62
1:D:111:HIS:C	1:D:113:GLU:H	2.03	0.62
1:D:165:PHE:CD2	1:D:165:PHE:C	2.73	0.62
1:D:158:LEU:HD23	1:D:163:ILE:HG21	1.82	0.61
1:B:243:ILE:N	1:B:243:ILE:HD13	2.14	0.61
1:D:205:PHE:CD1	1:D:259:PHE:CE1	2.88	0.61
1:C:168:THR:O	1:C:169:ASP:HB2	2.01	0.61
1:A:253:ASN:O	1:B:262:SER:HB3	2.01	0.61
1:B:273:LEU:HD23	1:B:275:ILE:HG13	1.82	0.61
1:D:185:PHE:CE2	1:D:281:MET:SD	2.94	0.61
1:B:97:PRO:HA	1:B:108:THR:HG22	1.81	0.61
1:C:156:THR:HG22	1:C:157:THR:N	2.15	0.61
1:D:135:CYS:HG	3:D:608:ZN:ZN	1.14	0.61
1:D:208:ILE:HG12	1:D:209:VAL:H	1.65	0.61
1:A:168:THR:HG22	2:E:122:PRO:HD3	1.83	0.61
1:A:178:TRP:N	1:A:178:TRP:CE3	2.69	0.61
1:B:236:TRP:CB	1:B:257:LEU:HB2	2.31	0.61
1:A:263:ILE:HG13	1:B:254:SER:HB3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:HIS:O	1:C:121:CYS:HB2	2.02	0.60
1:D:193:LEU:HD13	1:D:207:ALA:HB2	1.83	0.60
1:B:168:THR:CG2	2:F:120:VAL:O	2.43	0.60
1:B:214:THR:O	1:B:217:GLN:HB2	2.01	0.60
1:D:157:THR:HG22	1:D:158:LEU:N	2.15	0.59
1:D:184:CYS:O	1:D:185:PHE:HB2	2.02	0.59
1:A:176:VAL:CG1	2:E:122:PRO:HB3	2.33	0.59
1:A:257:LEU:C	1:A:258:VAL:CG1	2.71	0.59
1:B:224:ARG:HG3	1:B:237:GLU:HG3	1.84	0.59
1:B:263:ILE:O	1:B:265:GLN:N	2.36	0.58
1:C:236:TRP:HB3	1:C:257:LEU:HD13	1.85	0.58
1:A:257:LEU:C	1:A:258:VAL:HG12	2.23	0.58
1:A:178:TRP:HE3	1:A:178:TRP:H	1.51	0.58
1:C:128:CYS:C	1:C:130:CYS:H	2.06	0.58
1:C:165:PHE:CD2	1:C:165:PHE:C	2.77	0.58
1:B:263:ILE:C	1:B:265:GLN:N	2.58	0.57
1:D:158:LEU:HD23	1:D:163:ILE:HG23	1.86	0.57
1:D:185:PHE:HE2	1:D:281:MET:SD	2.27	0.57
1:C:101:ALA:HA	1:C:105:CYS:HB3	1.87	0.57
1:B:228:ASN:ND2	1:B:228:ASN:N	2.52	0.57
1:D:249:THR:O	1:D:253:ASN:ND2	2.38	0.57
1:C:231:ARG:NH1	1:D:239:THR:HG21	2.20	0.57
1:A:211:LEU:HD23	1:A:218:ALA:HA	1.86	0.57
1:D:214:THR:HG23	1:D:217:GLN:HG3	1.86	0.57
1:B:168:THR:HG23	2:F:122:PRO:HD3	1.87	0.56
1:D:259:PHE:HB2	1:D:263:ILE:HD12	1.86	0.56
1:C:210:GLN:NE2	1:C:243:ILE:HA	2.19	0.56
1:C:114:LYS:HE3	1:C:118:GLU:OE2	2.04	0.56
1:C:161:GLU:HA	1:C:182:GLN:NE2	2.20	0.56
1:D:194:GLU:OE2	1:D:196:GLN:NE2	2.38	0.56
1:D:265:GLN:HE21	1:D:265:GLN:H	1.53	0.56
1:A:256:CYS:SG	1:A:258:VAL:HG12	2.45	0.56
1:D:210:GLN:NE2	1:D:243:ILE:HG22	2.21	0.56
1:D:96:PHE:HE1	1:D:111:HIS:HB2	1.70	0.55
1:C:160:GLY:O	1:C:184:CYS:HB3	2.06	0.55
1:C:103:SER:HB2	1:C:123:PHE:HB3	1.89	0.55
1:C:206:PHE:CD2	1:C:251:ILE:HG23	2.42	0.55
1:C:234:LEU:HB2	1:D:236:TRP:HE3	1.72	0.55
1:B:184:CYS:O	1:B:185:PHE:HB2	2.06	0.55
1:B:260:ASP:O	1:B:263:ILE:HG22	2.07	0.55
1:D:214:THR:N	1:D:217:GLN:HE21	2.00	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:158:LEU:HG	1:D:163:ILE:HD13	1.89	0.55
1:D:96:PHE:CE1	1:D:111:HIS:HB2	2.42	0.55
1:A:95:LEU:HB2	1:A:108:THR:HG22	1.89	0.55
1:C:176:VAL:H	1:C:195:LYS:HB3	1.72	0.55
1:A:176:VAL:HG11	2:E:122:PRO:HB3	1.89	0.55
1:B:212:ILE:HA	1:B:243:ILE:HD12	1.87	0.55
1:B:168:THR:HA	1:B:272:ASN:OD1	2.07	0.54
1:D:208:ILE:HG12	1:D:209:VAL:N	2.23	0.54
1:D:232:ARG:NH1	1:D:232:ARG:HG3	2.19	0.54
1:B:244:HIS:O	1:B:246:GLY:N	2.40	0.54
1:B:260:ASP:N	1:B:260:ASP:OD1	2.41	0.53
1:A:242:SER:O	1:A:244:HIS:N	2.40	0.53
1:B:172:LEU:CD2	1:C:252:MET:HB3	2.38	0.53
1:B:124:ARG:HD2	1:B:125:PRO:O	2.08	0.53
1:A:247:ILE:O	1:A:251:ILE:HG13	2.08	0.53
1:A:269:GLU:C	1:A:270:ASN:HD22	2.12	0.53
1:C:165:PHE:HD2	1:C:165:PHE:C	2.11	0.53
1:D:169:ASP:HA	1:D:271:GLY:O	2.07	0.53
1:B:179:VAL:HG13	1:B:192:VAL:CG2	2.35	0.53
1:A:183:SER:O	1:A:184:CYS:HB3	2.09	0.52
1:B:92:ASN:C	1:B:94:VAL:H	2.13	0.52
1:C:247:ILE:O	1:C:251:ILE:HG13	2.09	0.52
1:D:111:HIS:O	1:D:114:LYS:HB2	2.09	0.52
1:A:276:ASN:N	1:A:276:ASN:HD22	2.07	0.52
1:C:183:SER:CB	1:C:188:HIS:ND1	2.72	0.52
1:B:236:TRP:CZ2	1:B:238:ALA:CB	2.91	0.52
1:D:210:GLN:HG2	1:D:247:ILE:HG21	1.91	0.52
1:A:171:ASN:O	1:A:172:LEU:C	2.46	0.52
1:A:158:LEU:HD21	2:E:116:PRO:HB2	1.91	0.52
1:D:141:LEU:HA	1:D:144:VAL:HG23	1.91	0.51
1:D:158:LEU:HD22	2:H:118:ALA:HB2	1.91	0.51
1:A:101:ALA:HA	1:A:105:CYS:H	1.76	0.51
1:C:145:MET:CE	1:C:212:ILE:HG13	2.41	0.51
1:D:168:THR:O	2:H:122:PRO:HG3	2.11	0.51
1:C:170:ILE:N	1:C:271:GLY:O	2.43	0.51
1:C:192:VAL:O	1:C:192:VAL:HG12	2.09	0.51
1:D:249:THR:O	1:D:252:MET:HB3	2.11	0.51
1:A:204:GLN:HB3	1:A:206:PHE:CE1	2.46	0.51
1:D:241:ARG:NH1	1:D:250:ALA:HB2	2.25	0.51
1:B:249:THR:O	1:B:253:ASN:ND2	2.42	0.51
1:D:172:LEU:HB3	1:D:176:VAL:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:266:LEU:N	1:D:266:LEU:HD22	2.27	0.50
1:D:241:ARG:CZ	1:D:250:ALA:HB2	2.41	0.50
1:A:259:PHE:HB2	1:A:263:ILE:CG2	2.39	0.50
1:B:253:ASN:ND2	1:B:253:ASN:H	2.09	0.50
1:A:168:THR:HB	2:E:120:VAL:O	2.12	0.50
1:A:221:PHE:O	1:A:239:THR:HG22	2.12	0.50
1:A:236:TRP:CZ2	1:A:238:ALA:HB2	2.47	0.50
1:A:263:ILE:O	1:A:264:ALA:C	2.49	0.50
1:C:145:MET:HE3	1:C:188:HIS:HB2	1.94	0.50
1:B:210:GLN:HE22	1:B:247:ILE:HG12	1.73	0.50
1:D:163:ILE:CG2	1:D:164:VAL:N	2.72	0.50
2:H:114:LEU:O	2:H:114:LEU:HD23	2.12	0.50
1:D:265:GLN:H	1:D:265:GLN:NE2	2.10	0.50
1:D:187:PHE:CD2	1:D:211:LEU:HD11	2.46	0.49
1:D:231:ARG:O	1:D:232:ARG:HG3	2.13	0.49
1:C:156:THR:CG2	1:C:157:THR:N	2.75	0.49
1:B:168:THR:CG2	2:F:122:PRO:HD3	2.43	0.49
1:D:276:ASN:OD1	2:H:115:ARG:NH1	2.46	0.49
1:A:119:GLU:HA	1:A:124:ARG:NH2	2.28	0.49
1:B:117:HIS:C	1:B:117:HIS:CD2	2.86	0.49
1:A:101:ALA:O	1:A:104:GLY:N	2.46	0.49
1:C:191:LEU:HD23	1:C:191:LEU:N	2.28	0.48
1:C:171:ASN:ND2	1:C:171:ASN:H	2.11	0.48
1:B:214:THR:OG1	1:B:217:GLN:HG3	2.13	0.48
1:C:270:ASN:O	1:C:272:ASN:N	2.46	0.48
1:C:231:ARG:NH1	1:D:219:GLU:OE1	2.46	0.48
1:A:145:MET:HG3	1:A:188:HIS:CD2	2.48	0.48
1:B:206:PHE:CD2	1:B:251:ILE:CG2	2.97	0.48
1:D:221:PHE:O	1:D:239:THR:HG22	2.13	0.48
1:A:143:ALA:C	1:A:146:PRO:HD2	2.33	0.48
1:D:157:THR:CG2	1:D:158:LEU:N	2.76	0.48
1:A:176:VAL:CG1	1:A:177:ASP:N	2.75	0.48
1:D:225:LEU:HD12	1:D:257:LEU:HB3	1.96	0.48
1:D:141:LEU:O	1:D:144:VAL:HG23	2.14	0.48
1:B:197:GLU:HA	1:B:203:GLN:HA	1.95	0.48
1:B:92:ASN:O	1:B:94:VAL:N	2.46	0.48
1:D:234:LEU:HD23	1:D:257:LEU:HD11	1.96	0.47
1:C:212:ILE:HA	1:C:243:ILE:CD1	2.26	0.47
1:B:269:GLU:O	1:B:272:ASN:HB2	2.14	0.47
1:B:253:ASN:HD22	1:B:253:ASN:H	1.62	0.47
1:A:218:ALA:HB1	1:A:240:PRO:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:LEU:HG	1:B:266:LEU:H	1.50	0.47
1:B:253:ASN:O	1:B:254:SER:CB	2.48	0.47
1:B:159:GLN:NE2	1:C:96:PHE:CE2	2.80	0.47
1:D:163:ILE:HG21	1:D:180:MET:HE1	1.97	0.47
1:A:101:ALA:HA	1:A:105:CYS:CB	2.42	0.47
1:B:273:LEU:HD23	1:B:275:ILE:CG1	2.43	0.47
1:A:101:ALA:CA	1:A:105:CYS:HB3	2.45	0.47
1:B:209:VAL:HG23	1:B:225:LEU:HD21	1.97	0.46
1:D:128:CYS:HA	1:D:129:PRO:HD3	1.83	0.46
1:D:168:THR:HG22	1:D:272:ASN:OD1	2.15	0.46
1:D:227:LEU:N	1:D:227:LEU:HD23	2.30	0.46
1:B:276:ASN:ND2	1:B:276:ASN:N	2.62	0.46
1:A:193:LEU:HD12	1:A:194:GLU:H	1.79	0.46
1:B:165:PHE:C	1:B:165:PHE:CD2	2.89	0.46
2:H:122:PRO:O	2:H:123:THR:OG1	2.30	0.46
1:A:103:SER:O	1:A:123:PHE:CE2	2.68	0.46
1:B:263:ILE:O	1:B:266:LEU:HG	2.15	0.46
1:B:247:ILE:CD1	1:B:251:ILE:HD11	2.44	0.46
1:C:107:ILE:N	1:C:107:ILE:CD1	2.63	0.46
1:A:209:VAL:HG12	1:A:210:GLN:N	2.31	0.46
1:A:232:ARG:HB2	1:B:236:TRP:HH2	1.79	0.46
1:C:265:GLN:HE21	1:C:265:GLN:N	2.14	0.45
1:D:209:VAL:HG23	1:D:225:LEU:CD2	2.43	0.45
1:D:188:HIS:HB2	1:D:212:ILE:HD12	1.98	0.45
1:B:263:ILE:HG13	1:B:266:LEU:HD11	1.98	0.45
1:B:250:ALA:O	1:B:254:SER:N	2.49	0.45
1:D:228:ASN:C	1:D:267:PHE:HD2	2.20	0.45
1:A:263:ILE:HA	1:A:266:LEU:HD12	1.98	0.45
1:B:128:CYS:SG	1:B:130:CYS:HB3	2.56	0.45
1:C:233:ARG:NH1	1:D:237:GLU:OE1	2.49	0.45
1:B:180:MET:HG3	1:B:181:MET:N	2.31	0.45
1:A:176:VAL:CG1	1:A:177:ASP:H	2.27	0.45
1:B:141:LEU:O	1:B:144:VAL:HG23	2.17	0.45
1:B:174:GLY:O	1:B:176:VAL:HG23	2.16	0.45
1:A:261:THR:O	1:A:265:GLN:HG2	2.16	0.45
1:B:227:LEU:CD2	1:B:275:ILE:HG12	2.47	0.45
1:B:165:PHE:HD2	1:B:165:PHE:C	2.19	0.45
1:A:145:MET:N	1:A:146:PRO:CD	2.79	0.45
1:D:214:THR:H	1:D:217:GLN:NE2	2.06	0.45
1:A:183:SER:OG	1:A:188:HIS:ND1	2.46	0.45
1:D:227:LEU:HD21	1:D:257:LEU:CD2	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:185:PHE:CD1	1:C:221:PHE:HE2	2.35	0.44
1:C:129:PRO:HG2	1:C:148:LEU:HD21	1.98	0.44
1:C:130:CYS:HA	1:C:131:PRO:HD2	1.81	0.44
1:C:239:THR:HB	1:C:240:PRO:HD2	1.99	0.44
1:A:202:HIS:HB2	1:B:199:TYR:CD2	2.53	0.44
1:D:263:ILE:HA	1:D:266:LEU:HD23	1.99	0.44
1:D:158:LEU:CD2	1:D:163:ILE:HG23	2.47	0.44
1:A:223:TYR:CE2	1:A:238:ALA:HB3	2.51	0.44
1:D:206:PHE:CB	1:D:251:ILE:HD12	2.45	0.44
1:D:263:ILE:O	1:D:266:LEU:HB2	2.17	0.44
1:C:185:PHE:O	1:C:187:PHE:CD1	2.69	0.44
1:C:259:PHE:HB2	1:C:263:ILE:HG21	2.00	0.44
1:D:128:CYS:SG	1:D:130:CYS:HB3	2.57	0.44
1:C:145:MET:HE2	1:C:212:ILE:HG13	2.00	0.43
1:C:166:LEU:HB3	2:G:119:MET:HA	1.99	0.43
1:C:241:ARG:HH11	1:C:241:ARG:HA	1.82	0.43
1:B:244:HIS:O	1:B:245:GLU:C	2.56	0.43
1:C:165:PHE:HA	1:C:180:MET:HE1	2.00	0.43
1:B:241:ARG:NH1	1:B:250:ALA:HB2	2.34	0.43
1:C:105:CYS:SG	1:C:107:ILE:HD11	2.59	0.43
1:C:195:LYS:HD2	1:C:205:PHE:CE2	2.54	0.43
1:C:236:TRP:CZ2	1:C:256:CYS:HA	2.53	0.43
1:A:111:HIS:CD2	2:H:116:PRO:HB3	2.53	0.43
1:A:140:SER:O	1:A:141:LEU:C	2.56	0.43
1:C:170:ILE:HA	1:C:170:ILE:HD12	1.82	0.43
1:D:101:ALA:HA	1:D:105:CYS:H	1.84	0.43
1:D:145:MET:N	1:D:146:PRO:HD2	2.33	0.42
1:B:163:ILE:HA	2:F:116:PRO:HD2	2.01	0.42
1:D:141:LEU:HD21	1:D:213:GLY:HA2	2.01	0.42
1:D:205:PHE:C	1:D:206:PHE:CD1	2.93	0.42
1:A:269:GLU:C	1:A:270:ASN:ND2	2.72	0.42
1:A:276:ASN:H	1:A:276:ASN:HD22	1.67	0.42
1:D:211:LEU:HD23	1:D:218:ALA:HB2	2.00	0.42
1:A:199:TYR:O	1:A:200:ASP:C	2.57	0.42
1:A:205:PHE:HE1	1:A:261:THR:HA	1.84	0.42
1:C:206:PHE:HD2	1:C:256:CYS:SG	2.42	0.42
1:A:197:GLU:HG2	1:A:197:GLU:H	1.76	0.42
1:D:206:PHE:CD2	1:D:251:ILE:HG23	2.54	0.42
1:D:111:HIS:C	1:D:113:GLU:N	2.72	0.42
1:B:171:ASN:O	1:B:172:LEU:C	2.57	0.42
1:B:126:TYR:O	1:B:138:GLN:HA	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:ALA:O	1:C:118:GLU:HB2	2.20	0.42
1:A:151:GLN:O	1:A:151:GLN:HG3	2.19	0.42
1:C:128:CYS:C	1:C:130:CYS:N	2.71	0.42
1:A:216:LYS:HD3	1:A:216:LYS:HA	1.82	0.42
1:B:169:ASP:HA	1:B:271:GLY:O	2.20	0.42
1:B:172:LEU:HA	1:B:173:PRO:HD3	1.90	0.42
1:D:182:GLN:HB3	1:D:191:LEU:HD21	2.02	0.42
1:B:224:ARG:HG3	1:B:237:GLU:CG	2.47	0.42
1:A:223:TYR:CZ	1:A:238:ALA:HB3	2.55	0.42
1:D:275:ILE:HG22	1:D:277:VAL:CG2	2.50	0.42
1:D:263:ILE:HG22	1:D:267:PHE:CE1	2.54	0.42
1:C:243:ILE:HG13	1:C:243:ILE:H	1.71	0.41
1:A:263:ILE:HG22	1:A:264:ALA:N	2.34	0.41
1:A:257:LEU:HG	1:A:258:VAL:N	2.35	0.41
1:D:265:GLN:N	1:D:265:GLN:NE2	2.68	0.41
1:D:234:LEU:CD2	1:D:257:LEU:HD11	2.49	0.41
1:C:168:THR:O	1:C:169:ASP:CB	2.68	0.41
1:C:185:PHE:O	1:C:186:GLY:C	2.59	0.41
1:A:185:PHE:O	1:A:187:PHE:CD1	2.73	0.41
1:B:208:ILE:HG12	1:B:209:VAL:N	2.34	0.41
1:C:166:LEU:HD23	2:G:119:MET:HE1	2.02	0.41
1:A:165:PHE:C	1:A:165:PHE:CD2	2.93	0.41
1:B:206:PHE:CD2	1:B:251:ILE:HG23	2.56	0.41
1:A:157:THR:HG22	1:A:158:LEU:N	2.34	0.41
1:A:103:SER:O	1:A:123:PHE:CD2	2.73	0.41
1:A:257:LEU:O	1:A:258:VAL:HG12	2.20	0.41
1:A:246:GLY:C	1:A:247:ILE:HG23	2.41	0.41
1:B:265:GLN:OE1	1:B:265:GLN:HA	2.19	0.41
1:A:211:LEU:CD2	1:A:218:ALA:HA	2.49	0.41
1:B:169:ASP:HB3	1:C:253:ASN:HD21	1.85	0.41
1:A:260:ASP:O	1:A:263:ILE:HB	2.20	0.41
1:D:206:PHE:CD2	1:D:251:ILE:HD12	2.56	0.41
1:D:266:LEU:N	1:D:266:LEU:CD2	2.84	0.41
1:D:157:THR:HA	1:D:181:MET:HB3	2.03	0.41
1:C:176:VAL:O	1:C:177:ASP:OD1	2.39	0.41
1:C:263:ILE:HG13	1:D:254:SER:O	2.21	0.41
1:C:172:LEU:HD23	1:C:173:PRO:HD2	2.02	0.41
1:D:145:MET:CE	1:D:145:MET:CA	2.95	0.41
1:D:180:MET:CG	1:D:181:MET:N	2.81	0.41
1:C:193:LEU:HD12	1:C:194:GLU:H	1.78	0.41
1:A:270:ASN:N	1:A:270:ASN:ND2	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:169:ASP:CA	1:D:271:GLY:O	2.69	0.41
1:A:191:LEU:N	1:A:191:LEU:HD23	2.35	0.41
1:B:146:PRO:O	1:B:150:HIS:CD2	2.74	0.41
1:B:255:ASP:O	1:B:256:CYS:HB3	2.21	0.41
1:B:249:THR:HG23	1:B:250:ALA:N	2.36	0.40
1:B:261:THR:O	1:B:265:GLN:HG2	2.21	0.40
1:A:208:ILE:HD12	1:A:256:CYS:HB3	2.03	0.40
1:A:209:VAL:CG1	1:A:210:GLN:N	2.84	0.40
1:A:170:ILE:HA	1:A:170:ILE:HD13	1.70	0.40
1:D:167:ALA:HB1	1:D:178:TRP:CD1	2.56	0.40
1:B:276:ASN:HB3	2:F:115:ARG:NH1	2.35	0.40
1:A:276:ASN:N	1:A:276:ASN:ND2	2.68	0.40
1:C:128:CYS:SG	1:C:130:CYS:HB2	2.61	0.40
1:C:234:LEU:HB2	1:D:236:TRP:CE3	2.52	0.40
1:A:210:GLN:HE21	1:A:210:GLN:HB3	1.58	0.40
1:D:141:LEU:HA	1:D:144:VAL:CG2	2.51	0.40
1:C:171:ASN:ND2	1:C:171:ASN:N	2.70	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	189/191 (99%)	159 (84%)	25 (13%)	5 (3%)	7	33
1	B	189/191 (99%)	160 (85%)	22 (12%)	7 (4%)	4	23
1	C	189/191 (99%)	157 (83%)	25 (13%)	7 (4%)	4	23
1	D	189/191 (99%)	153 (81%)	29 (15%)	7 (4%)	4	23
2	E	9/24 (38%)	8 (89%)	1 (11%)	0	100	100
2	F	9/24 (38%)	7 (78%)	2 (22%)	0	100	100
2	G	9/24 (38%)	8 (89%)	1 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	9/24 (38%)	7 (78%)	0	2 (22%)	0	0
All	All	792/860 (92%)	659 (83%)	105 (13%)	28 (4%)	4	24

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	ILE
1	B	244	HIS
1	B	245	GLU
1	B	256	CYS
1	C	155	ILE
1	A	245	GLU
1	B	93	SER
1	B	264	ALA
1	C	175	ALA
1	D	131	PRO
1	D	132	GLY
1	D	256	CYS
1	D	261	THR
1	B	200	ASP
1	C	161	GLU
1	C	198	LYS
2	H	122	PRO
2	H	123	THR
1	A	98	CYS
1	A	133	ALA
1	B	270	ASN
1	D	112	THR
1	D	238	ALA
1	A	247	ILE
1	C	170	ILE
1	C	186	GLY
1	C	241	ARG
1	D	212	ILE

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	165/165 (100%)	132 (80%)	33 (20%)	1	8
1	B	165/165 (100%)	135 (82%)	30 (18%)	2	11
1	C	165/165 (100%)	141 (86%)	24 (14%)	4	18
1	D	165/165 (100%)	133 (81%)	32 (19%)	2	9
2	E	10/23 (44%)	8 (80%)	2 (20%)	1	8
2	F	10/23 (44%)	9 (90%)	1 (10%)	9	34
2	G	10/23 (44%)	8 (80%)	2 (20%)	1	8
2	H	10/23 (44%)	7 (70%)	3 (30%)	0	2
All	All	700/752 (93%)	573 (82%)	127 (18%)	2	11

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	SER
1	A	95	LEU
1	A	99	LYS
1	A	103	SER
1	A	108	THR
1	A	112	THR
1	A	120	LEU
1	A	130	CYS
1	A	134	SER
1	A	141	LEU
1	A	145	MET
1	A	150	HIS
1	A	154	SER
1	A	165	PHE
1	A	168	THR
1	A	170	ILE
1	A	181	MET
1	A	191	LEU
1	A	199	TYR
1	A	210	GLN
1	A	214	THR
1	A	215	ARG
1	A	219	GLU
1	A	232	ARG
1	A	233	ARG

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Mol	Chain	Res	Type
1	A	234	LEU
1	A	237	GLU
1	A	240	PRO
1	A	254	SER
1	A	257	LEU
1	A	261	THR
1	A	270	ASN
1	A	282	CYS
1	B	94	VAL
1	B	109	LEU
1	B	111	HIS
1	B	145	MET
1	B	152	HIS
1	B	154	SER
1	B	156	THR
1	B	159	GLN
1	B	165	PHE
1	B	177	ASP
1	B	179	VAL
1	B	180	MET
1	B	183	SER
1	B	190	MET
1	B	191	LEU
1	B	214	THR
1	B	215	ARG
1	B	228	ASN
1	B	239	THR
1	B	243	ILE
1	B	244	HIS
1	B	245	GLU
1	B	251	ILE
1	B	253	ASN
1	B	254	SER
1	B	260	ASP
1	B	261	THR
1	B	262	SER
1	B	266	LEU
1	B	276	ASN
1	C	98	CYS
1	C	100	TYR
1	C	107	ILE
1	C	114	LYS

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Mol	Chain	Res	Type
1	C	141	LEU
1	C	153	LYS
1	C	161	GLU
1	C	165	PHE
1	C	169	ASP
1	C	170	ILE
1	C	171	ASN
1	C	172	LEU
1	C	183	SER
1	C	191	LEU
1	C	193	LEU
1	C	199	TYR
1	C	211	LEU
1	C	216	LYS
1	C	235	THR
1	C	236	TRP
1	C	241	ARG
1	C	260	ASP
1	C	263	ILE
1	C	281	MET
1	D	93	SER
1	D	103	SER
1	D	128	CYS
1	D	140	SER
1	D	145	MET
1	D	149	MET
1	D	152	HIS
1	D	156	THR
1	D	165	PHE
1	D	168	THR
1	D	169	ASP
1	D	170	ILE
1	D	178	TRP
1	D	180	MET
1	D	181	MET
1	D	183	SER
1	D	191	LEU
1	D	202	HIS
1	D	203	GLN
1	D	208	ILE
1	D	210	GLN
1	D	214	THR

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Mol	Chain	Res	Type
1	D	227	LEU
1	D	232	ARG
1	D	237	GLU
1	D	239	THR
1	D	249	THR
1	D	253	ASN
1	D	261	THR
1	D	263	ILE
1	D	265	GLN
1	D	270	ASN
2	E	115	ARG
2	E	124	VAL
2	F	114	LEU
2	G	119	MET
2	G	123	THR
2	H	114	LEU
2	H	115	ARG
2	H	117	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	GLN
1	A	253	ASN
1	A	270	ASN
1	A	276	ASN
1	B	111	HIS
1	B	159	GLN
1	B	196	GLN
1	B	228	ASN
1	B	253	ASN
1	B	270	ASN
1	C	171	ASN
1	C	182	GLN
1	C	228	ASN
1	C	253	ASN
1	C	265	GLN
1	D	152	HIS
1	D	171	ASN
1	D	182	GLN
1	D	196	GLN
1	D	210	GLN

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Mol	Chain	Res	Type
1	D	217	GLN
1	D	253	ASN
1	D	265	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	191/191 (100%)	-0.03	0 100 100	13, 66, 99, 133	0
1	B	191/191 (100%)	-0.06	1 (0%) 91 76	14, 71, 112, 126	0
1	C	191/191 (100%)	0.01	3 (1%) 74 47	14, 72, 109, 148	0
1	D	191/191 (100%)	0.04	2 (1%) 84 60	25, 83, 137, 143	0
2	E	11/24 (45%)	0.30	0 100 100	48, 56, 69, 75	0
2	F	11/24 (45%)	0.23	0 100 100	42, 68, 92, 93	0
2	G	11/24 (45%)	0.30	0 100 100	52, 71, 94, 101	0
2	H	11/24 (45%)	0.25	1 (9%) 11 4	65, 77, 95, 107	0
All	All	808/860 (93%)	0.00	7 (0%) 85 64	13, 72, 126, 148	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	201	GLY	3.9
2	H	124	VAL	3.0
1	B	198	LYS	2.6
1	C	202	HIS	2.4
1	D	126	TYR	2.3
1	D	199	TYR	2.1
1	C	200	ASP	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	ZN	A	602	1/1	0.98	0.21	1.01	78,78,78,78	0
3	ZN	C	605	1/1	0.99	0.18	0.31	78,78,78,78	0
3	ZN	B	604	1/1	0.98	0.21	0.13	97,97,97,97	0
3	ZN	D	607	1/1	0.91	0.18	-0.26	144,144,144,144	0
3	ZN	C	606	1/1	0.97	0.14	-0.73	95,95,95,95	0
3	ZN	D	608	1/1	0.98	0.14	-0.95	115,115,115,115	0
3	ZN	B	603	1/1	0.99	0.18	-	89,89,89,89	0
3	ZN	A	601	1/1	0.98	0.20	-	80,80,80,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.