



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 15, 2023 – 01:00 AM JST

PDB ID : 6IQD
Title : Crystal structure of Alcohol dehydrogenase from *Geobacillus stearothermophilus*
Authors : Xue, S.; Feng, Y.; Guo, X.; Zhao, Z.
Deposited on : 2018-11-07
Resolution : 2.84 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

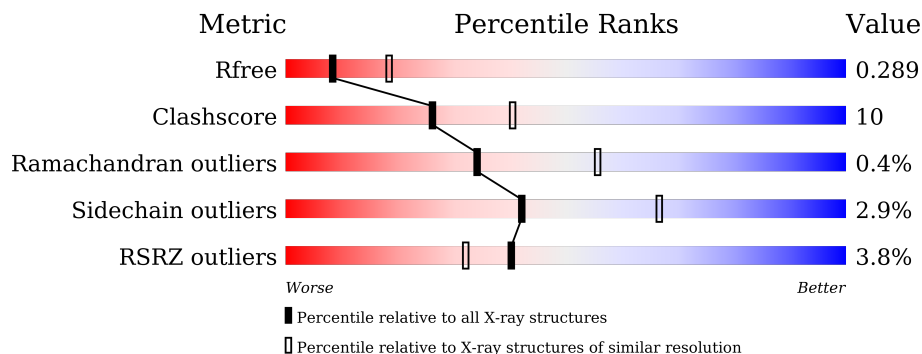
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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.



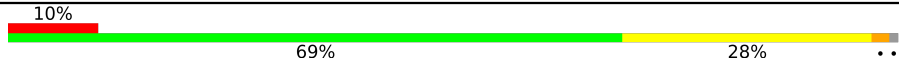

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1031 (2.86-2.82)
Clashscore	141614	1078 (2.86-2.82)
Ramachandran outliers	138981	1050 (2.86-2.82)
Sidechain outliers	138945	1051 (2.86-2.82)
RSRZ outliers	127900	1019 (2.86-2.82)

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	339	 83% 12% . .
1	B	339	 81% 18% ..
1	C	339	 3% 79% 19% ..
1	D	339	 2% 83% 15% .
1	E	339	 73% 23% ..
1	F	339	 2% 81% 17% ..

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Mol	Chain	Length	Quality of chain
1	G	339	
1	H	339	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 20152 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 Alcohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	336	2517	1606	429	470	12	0	0	0
1	B	336	2517	1606	429	470	12	0	0	0
1	C	336	2517	1606	429	470	12	0	0	0
1	D	336	2517	1606	429	470	12	0	0	0
1	E	336	2517	1606	429	470	12	0	0	0
1	F	336	2517	1606	429	470	12	0	0	0
1	G	336	2517	1606	429	470	12	0	0	0
1	H	336	2517	1606	429	470	12	0	0	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

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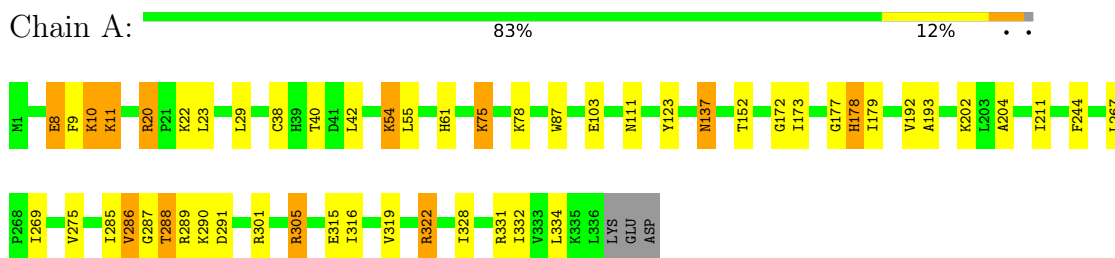
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total 2	Zn 2	0	0
2	H	2	Total 2	Zn 2	0	0

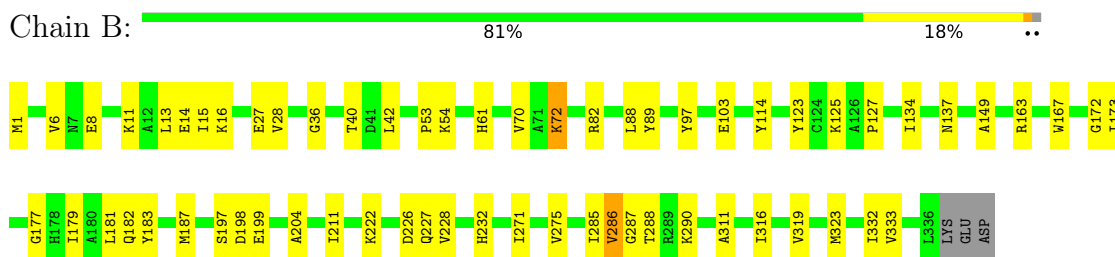
3 Residue-property plots [i](#)

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.

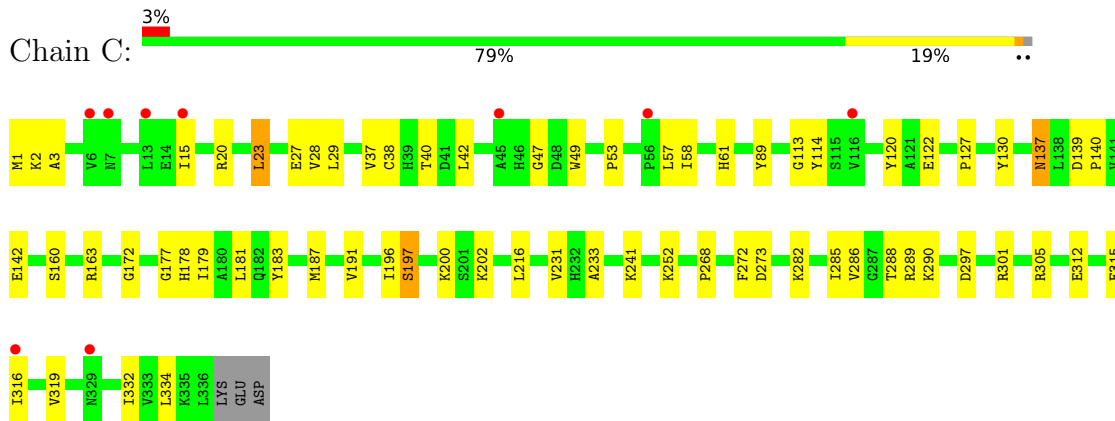
- Molecule 1: Alcohol dehydrogenase



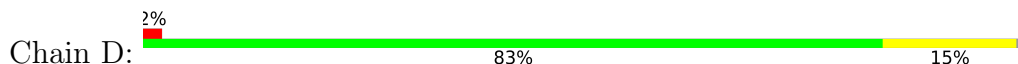
- Molecule 1: Alcohol dehydrogenase

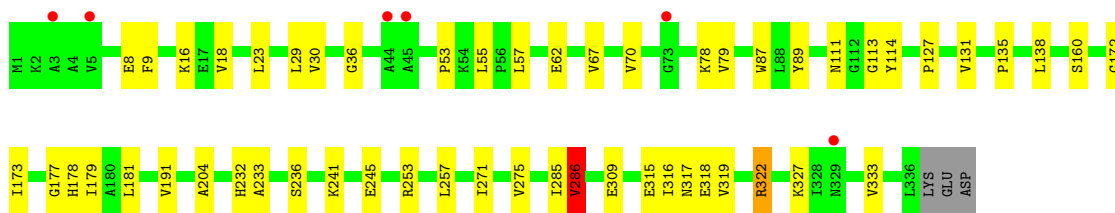


- Molecule 1: Alcohol dehydrogenase



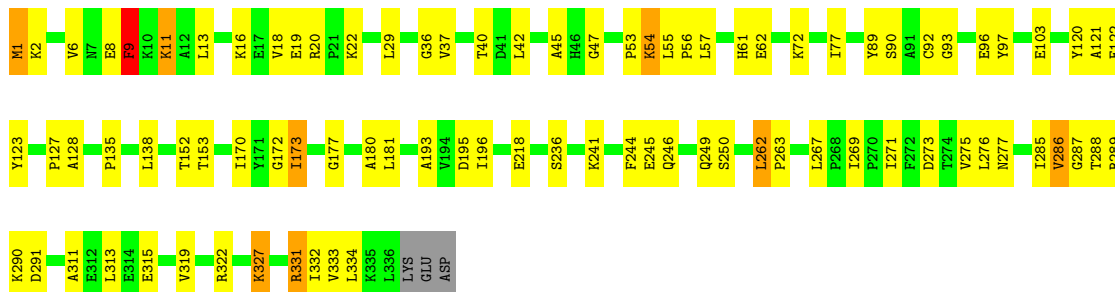
- Molecule 1: Alcohol dehydrogenase





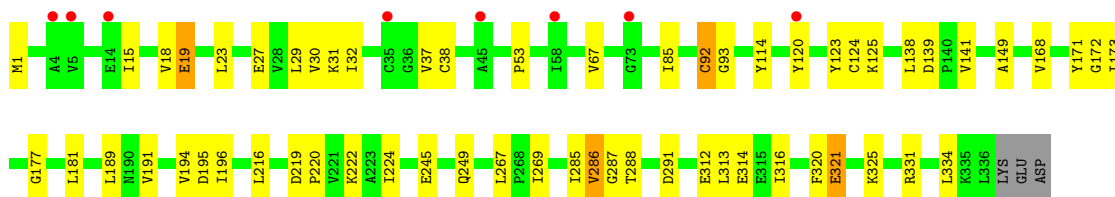
- Molecule 1: Alcohol dehydrogenase

Chain E: 73% 23% ..



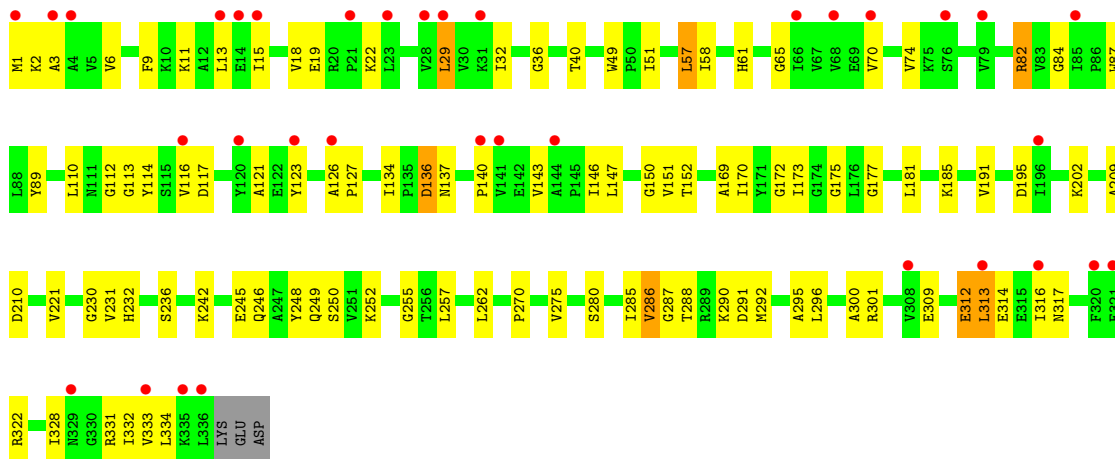
- Molecule 1: Alcohol dehydrogenase

Chain F: 81% 17% ..




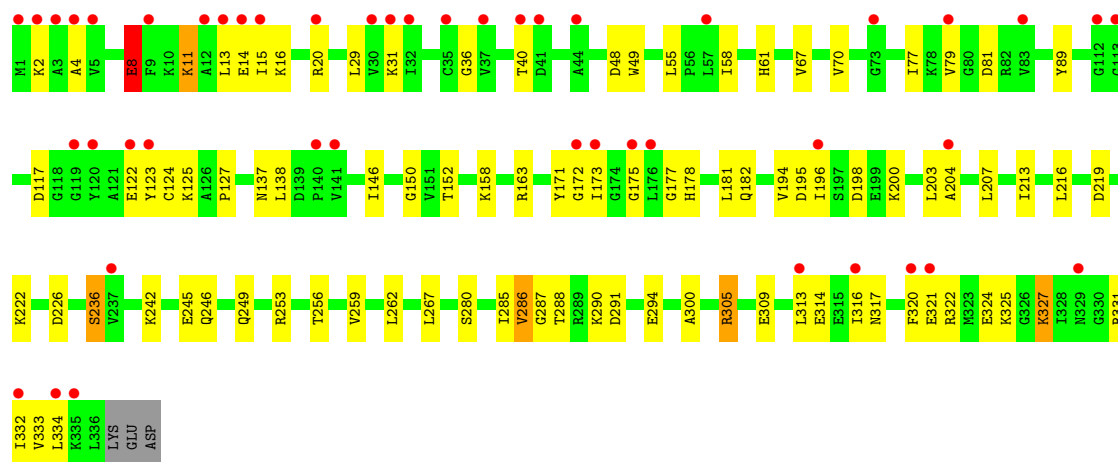
- Molecule 1: Alcohol dehydrogenase

Chain G: 69% 28% ..



- Molecule 1: Alcohol dehydrogenase

Chain H:  14% 72% 26% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.44Å 149.19Å 269.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.80 – 2.84 49.98 – 2.84	Depositor EDS
% Data completeness (in resolution range)	97.7 (49.80-2.84) 97.8 (49.98-2.84)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.86Å)	Xtrriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.232 , 0.291 0.232 , 0.289	Depositor DCC
R_{free} test set	3426 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtrriage
Anisotropy	0.034	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	20152	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% 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

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	0.33	0/2562	0.50	0/3471
1	B	0.30	0/2562	0.47	0/3471
1	C	0.29	0/2562	0.48	0/3471
1	D	0.29	0/2562	0.47	0/3471
1	E	0.30	0/2562	0.52	2/3471 (0.1%)
1	F	0.27	0/2562	0.49	1/3471 (0.0%)
1	G	0.28	0/2562	0.51	1/3471 (0.0%)
1	H	0.29	0/2562	0.53	1/3471 (0.0%)
All	All	0.29	0/20496	0.50	5/27768 (0.0%)

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	A	0	3

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	8	GLU	CB-CA-C	5.48	121.36	110.40
1	E	9	PHE	CB-CA-C	5.40	121.20	110.40
1	F	138	LEU	C-N-CA	5.17	134.62	121.70
1	G	57	LEU	CA-CB-CG	5.16	127.18	115.30
1	E	262	LEU	CA-CB-CG	-5.00	103.79	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	20	ARG	Sidechain
1	A	301	ARG	Sidechain
1	A	305	ARG	Sidechain

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	2517	0	2582	39	0
1	B	2517	0	2582	37	0
1	C	2517	0	2580	41	0
1	D	2517	0	2582	35	0
1	E	2517	0	2582	86	0
1	F	2517	0	2582	40	0
1	G	2517	0	2582	88	0
1	H	2517	0	2582	72	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
All	All	20152	0	20654	425	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 10.

All (425) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:9:PHE:CE1	1:E:55:LEU:HD21	1.39	1.52
1:E:9:PHE:HE1	1:E:55:LEU:CD2	1.42	1.33
1:E:9:PHE:CZ	1:E:55:LEU:HD21	1.65	1.30
1:E:9:PHE:CE1	1:E:55:LEU:CD2	2.14	1.26
1:E:9:PHE:HE1	1:E:55:LEU:CG	1.65	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:328:ILE:HD13	1:G:332:ILE:HD13	1.39	1.03
1:H:321:GLU:HA	1:H:324:GLU:OE2	1.59	0.99
1:G:312:GLU:OE2	1:G:314:GLU:OE1	1.82	0.95
1:E:9:PHE:CE1	1:E:55:LEU:CG	2.45	0.94
1:H:321:GLU:HG3	1:H:324:GLU:OE2	1.68	0.94
1:A:173:ILE:HD11	1:A:193:ALA:HB1	1.51	0.92
1:H:321:GLU:CA	1:H:324:GLU:OE2	2.18	0.91
1:G:146:ILE:HD12	1:G:150:GLY:HA3	1.53	0.90
1:G:312:GLU:OE1	1:G:313:LEU:N	2.06	0.89
1:G:328:ILE:HD13	1:G:332:ILE:CD1	2.04	0.88
1:E:153:THR:HG21	1:E:180:ALA:HA	1.55	0.87
1:G:6:VAL:HG12	1:G:13:LEU:HG	1.58	0.84
1:G:328:ILE:CD1	1:G:332:ILE:HD13	2.09	0.83
1:A:103:GLU:OE2	1:A:289:ARG:NE	2.11	0.82
1:A:11:LYS:HD2	1:A:11:LYS:N	1.96	0.80
1:E:9:PHE:CZ	1:E:55:LEU:CD2	2.56	0.80
1:G:32:ILE:HA	1:G:65:GLY:HA3	1.65	0.79
1:H:288:THR:HG22	1:H:291:ASP:CG	2.03	0.79
1:F:312:GLU:HG2	1:F:313:LEU:H	1.51	0.76
1:E:9:PHE:HD2	1:E:47:GLY:HA3	1.49	0.76
1:F:23:LEU:HD21	1:F:29:LEU:HB2	1.69	0.75
1:G:185:LYS:NZ	1:G:210:ASP:OD2	2.17	0.74
1:A:288:THR:HG22	1:A:291:ASP:H	1.53	0.74
1:D:285:ILE:HG23	1:D:286:VAL:HG23	1.67	0.74
1:D:36:GLY:HA2	1:D:333:VAL:HG12	1.70	0.73
1:E:9:PHE:HE1	1:E:55:LEU:HG	1.54	0.73
1:A:23:LEU:HD21	1:A:29:LEU:HB2	1.73	0.70
1:F:32:ILE:HD11	1:F:124:CYS:HB2	1.73	0.70
1:E:103:GLU:OE2	1:E:288:THR:OG1	2.09	0.70
1:B:173:ILE:HD13	1:B:204:ALA:HB2	1.73	0.69
1:H:15:ILE:HG12	1:H:314:GLU:HG3	1.74	0.69
1:G:232:HIS:CD2	1:G:255:GLY:HA2	2.28	0.69
1:E:153:THR:CG2	1:E:180:ALA:HA	2.20	0.69
1:A:8:GLU:HA	1:A:55:LEU:HD13	1.75	0.69
1:E:103:GLU:OE1	1:E:289:ARG:NE	2.25	0.69
1:E:9:PHE:CE1	1:E:55:LEU:HG	2.27	0.69
1:H:146:ILE:HG23	1:H:150:GLY:HA3	1.74	0.69
1:E:9:PHE:CE1	1:E:55:LEU:HD11	2.28	0.69
1:E:20:ARG:HG3	1:E:22:LYS:NZ	2.08	0.69
1:H:182:GLN:OE1	1:H:305:ARG:N	2.21	0.68
1:E:275:VAL:HG23	1:E:276:LEU:HD12	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:62:GLU:OE2	1:E:331:ARG:NH2	2.28	0.67
1:D:67:VAL:HG23	1:D:79:VAL:HA	1.76	0.67
1:A:22:LYS:HD2	1:A:23:LEU:H	1.58	0.67
1:D:8:GLU:HA	1:D:55:LEU:HD13	1.75	0.67
1:G:175:GLY:HA3	1:G:331:ARG:HE	1.59	0.67
1:F:85:ILE:HD12	1:F:85:ILE:O	1.94	0.67
1:G:173:ILE:HG22	1:G:195:ASP:HB2	1.77	0.67
1:A:20:ARG:CG	1:A:20:ARG:HH11	2.08	0.66
1:C:316:ILE:HD13	1:C:334:LEU:HD11	1.76	0.66
1:D:23:LEU:HD21	1:D:29:LEU:HB2	1.77	0.65
1:E:9:PHE:CE1	1:E:55:LEU:CD1	2.80	0.65
1:G:19:GLU:OE2	1:G:19:GLU:N	2.29	0.65
1:F:316:ILE:HD13	1:F:334:LEU:HD11	1.78	0.65
1:G:312:GLU:CD	1:G:313:LEU:H	1.97	0.65
1:A:173:ILE:HD13	1:A:204:ALA:HB2	1.78	0.65
1:B:288:THR:HG22	1:B:290:LYS:H	1.62	0.65
1:E:9:PHE:CD2	1:E:47:GLY:HA3	2.30	0.65
1:H:321:GLU:HG3	1:H:324:GLU:CD	2.18	0.64
1:F:312:GLU:OE2	1:F:314:GLU:OE2	2.16	0.64
1:C:23:LEU:HD23	1:C:27:GLU:HB2	1.77	0.64
1:G:15:ILE:HG12	1:G:314:GLU:HA	1.80	0.64
1:A:40:THR:HB	1:A:61:HIS:HE1	1.63	0.64
1:A:173:ILE:CD1	1:A:193:ALA:HB1	2.26	0.63
1:H:288:THR:HG22	1:H:291:ASP:OD2	1.98	0.63
1:G:1:MET:O	1:G:18:VAL:N	2.29	0.63
1:C:288:THR:HG22	1:C:290:LYS:H	1.65	0.62
1:E:173:ILE:HD11	1:E:193:ALA:HB1	1.80	0.62
1:F:312:GLU:HG2	1:F:313:LEU:N	2.15	0.62
1:G:316:ILE:HD13	1:G:334:LEU:HD11	1.82	0.62
1:C:23:LEU:HD21	1:C:27:GLU:C	2.20	0.61
1:G:15:ILE:HG13	1:G:313:LEU:HD23	1.81	0.61
1:C:1:MET:HG3	1:C:122:GLU:HB2	1.82	0.61
1:B:40:THR:HB	1:B:61:HIS:HE1	1.66	0.61
1:C:23:LEU:CD2	1:C:27:GLU:HB2	2.29	0.61
1:E:322:ARG:HG3	1:E:327:LYS:NZ	2.15	0.61
1:G:6:VAL:HG22	1:G:57:LEU:HB3	1.81	0.61
1:H:194:VAL:HG22	1:H:213:ILE:HB	1.80	0.61
1:B:72:LYS:HG3	1:B:72:LYS:O	2.01	0.61
1:H:288:THR:HG23	1:H:291:ASP:H	1.66	0.61
1:E:9:PHE:O	1:E:45:ALA:O	2.18	0.60
1:E:172:GLY:O	1:E:177:GLY:HA3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ARG:HH11	1:A:20:ARG:HG3	1.66	0.60
1:E:6:VAL:HA	1:E:13:LEU:HD22	1.83	0.60
1:G:40:THR:HB	1:G:61:HIS:HE1	1.66	0.60
1:C:40:THR:HG22	1:C:49:TRP:HH2	1.66	0.60
1:A:29:LEU:HD11	1:A:123:TYR:HB3	1.84	0.59
1:E:322:ARG:HG3	1:E:327:LYS:HD2	1.84	0.59
1:C:2:LYS:HG3	1:C:15:ILE:HD11	1.84	0.59
1:C:20:ARG:NH1	1:C:29:LEU:CD1	2.66	0.59
1:E:54:LYS:HD2	1:E:55:LEU:N	2.18	0.59
1:H:321:GLU:CG	1:H:324:GLU:OE2	2.47	0.59
1:C:241:LYS:HE2	1:C:268:PRO:HG2	1.85	0.58
1:E:322:ARG:HG3	1:E:327:LYS:CD	2.32	0.58
1:B:14:GLU:HG2	1:B:16:LYS:HB3	1.85	0.58
1:D:318:GLU:O	1:D:322:ARG:HB2	2.03	0.58
1:G:36:GLY:HA2	1:G:333:VAL:HG22	1.85	0.58
1:H:316:ILE:HD13	1:H:334:LEU:HD11	1.86	0.58
1:C:20:ARG:NH1	1:C:29:LEU:HD11	2.18	0.58
1:E:9:PHE:CD1	1:E:9:PHE:N	2.72	0.58
1:G:1:MET:SD	1:G:123:TYR:HB2	2.44	0.58
1:C:2:LYS:HB3	1:C:122:GLU:HG3	1.86	0.57
1:C:181:LEU:HD23	1:C:191:VAL:HG11	1.86	0.57
1:E:262:LEU:HD11	1:E:285:ILE:HD13	1.87	0.57
1:H:171:TYR:CE2	1:H:246:GLN:HB3	2.39	0.57
1:C:3:ALA:HB1	1:C:58:ILE:HD12	1.86	0.57
1:H:321:GLU:N	1:H:324:GLU:OE2	2.38	0.57
1:G:328:ILE:CD1	1:G:332:ILE:CD1	2.76	0.56
1:D:181:LEU:HD23	1:D:191:VAL:HG11	1.88	0.56
1:D:241:LYS:O	1:D:245:GLU:HG2	2.05	0.56
1:F:37:VAL:HG12	1:F:120:TYR:OH	2.06	0.56
1:E:9:PHE:N	1:E:9:PHE:HD1	2.04	0.56
1:A:269:ILE:HD12	1:F:269:ILE:HD12	1.88	0.55
1:F:286:VAL:HG12	1:F:287:GLY:N	2.21	0.55
1:C:288:THR:HG22	1:C:290:LYS:N	2.22	0.55
1:E:263:PRO:HG3	1:E:267:LEU:HD11	1.89	0.55
1:F:286:VAL:HG12	1:F:287:GLY:H	1.71	0.55
1:H:309:GLU:HB3	1:H:332:ILE:HD13	1.86	0.55
1:E:77:ILE:HD12	1:E:128:ALA:HB1	1.88	0.55
1:G:152:THR:HG23	1:G:285:ILE:HA	1.88	0.55
1:H:173:ILE:HG22	1:H:195:ASP:HB2	1.89	0.55
1:G:312:GLU:OE2	1:G:314:GLU:CD	2.44	0.55
1:A:54:LYS:HG3	1:A:54:LYS:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:20:ARG:HG3	1:E:22:LYS:HZ1	1.71	0.54
1:B:311:ALA:HB2	1:B:332:ILE:HD11	1.90	0.54
1:D:173:ILE:HD13	1:D:204:ALA:HB2	1.87	0.54
1:G:312:GLU:CD	1:G:313:LEU:N	2.57	0.54
1:C:27:GLU:HA	1:C:127:PRO:HA	1.89	0.54
1:D:322:ARG:HG2	1:D:327:LYS:HG2	1.90	0.54
1:G:181:LEU:HD23	1:G:191:VAL:HG11	1.90	0.54
1:H:13:LEU:N	1:H:317:ASN:HD21	2.06	0.54
1:G:51:ILE:O	1:G:114:TYR:OH	2.18	0.54
1:A:172:GLY:O	1:A:177:GLY:HA3	2.08	0.53
1:E:9:PHE:HD1	1:E:9:PHE:H	1.57	0.53
1:F:286:VAL:CG1	1:F:287:GLY:H	2.22	0.53
1:E:54:LYS:HD2	1:E:55:LEU:H	1.73	0.53
1:G:58:ILE:HG12	1:G:117:ASP:HB3	1.91	0.53
1:E:275:VAL:HG11	1:H:259:VAL:HG12	1.91	0.53
1:E:245:GLU:O	1:E:249:GLN:HG3	2.08	0.53
1:C:172:GLY:O	1:C:177:GLY:HA3	2.09	0.52
1:H:181:LEU:HD11	1:H:207:LEU:HB2	1.90	0.52
1:E:8:GLU:HB2	1:E:11:LYS:HG3	1.91	0.52
1:G:143:VAL:O	1:G:146:ILE:HG22	2.08	0.52
1:G:151:VAL:HG12	1:G:295:ALA:HB3	1.92	0.52
1:G:312:GLU:OE1	1:G:313:LEU:CB	2.58	0.52
1:A:173:ILE:O	1:A:173:ILE:HG22	2.09	0.52
1:A:20:ARG:CG	1:A:20:ARG:NH1	2.72	0.52
1:G:232:HIS:HD2	1:G:255:GLY:HA2	1.75	0.52
1:B:82:ARG:HG2	1:B:134:ILE:HB	1.91	0.52
1:B:53:PRO:HD3	1:B:114:TYR:CZ	2.45	0.52
1:G:242:LYS:O	1:G:246:GLN:HG3	2.10	0.52
1:H:31:LYS:HE3	1:H:123:TYR:OH	2.10	0.52
1:B:40:THR:HB	1:B:61:HIS:CE1	2.46	0.51
1:E:20:ARG:HG3	1:E:22:LYS:HZ3	1.75	0.51
1:E:36:GLY:HA2	1:E:333:VAL:HA	1.92	0.51
1:E:271:ILE:HD13	1:H:267:LEU:HG	1.93	0.51
1:B:211:ILE:HD13	1:B:227:GLN:NE2	2.26	0.51
1:E:288:THR:HG22	1:E:291:ASP:CG	2.31	0.51
1:B:275:VAL:HG12	1:D:285:ILE:HB	1.93	0.51
1:E:9:PHE:HZ	1:E:55:LEU:HD21	1.57	0.51
1:G:112:GLY:HA2	1:G:116:VAL:HG12	1.93	0.51
1:A:11:LYS:N	1:A:11:LYS:CD	2.72	0.50
1:F:15:ILE:HG12	1:F:314:GLU:HA	1.93	0.50
1:G:89:TYR:CZ	1:G:127:PRO:HD3	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:SER:HB3	1:C:200:LYS:HG3	1.93	0.50
1:C:285:ILE:HB	1:G:275:VAL:HG12	1.93	0.50
1:D:30:VAL:HG22	1:D:67:VAL:HG12	1.94	0.50
1:F:29:LEU:HD11	1:F:123:TYR:HB3	1.94	0.50
1:G:172:GLY:O	1:G:177:GLY:HA3	2.12	0.50
1:B:167:TRP:CG	1:B:228:VAL:HG21	2.47	0.50
1:G:57:LEU:HD21	1:G:113:GLY:O	2.12	0.50
1:H:2:LYS:HB2	1:H:122:GLU:HG3	1.94	0.50
1:E:16:LYS:HE3	1:E:18:VAL:HG22	1.93	0.49
1:G:170:ILE:HD13	1:G:191:VAL:HG13	1.94	0.49
1:H:31:LYS:HE3	1:H:123:TYR:CZ	2.47	0.49
1:H:8:GLU:HA	1:H:55:LEU:HD13	1.94	0.49
1:H:11:LYS:N	1:H:11:LYS:CD	2.74	0.49
1:B:319:VAL:CG2	1:B:332:ILE:HG21	2.43	0.49
1:D:16:LYS:HE2	1:D:18:VAL:HG12	1.94	0.49
1:E:288:THR:HG23	1:E:290:LYS:H	1.77	0.49
1:F:1:MET:SD	1:F:123:TYR:HB2	2.53	0.49
1:H:36:GLY:HA2	1:H:333:VAL:HA	1.94	0.49
1:A:8:GLU:OE1	1:A:9:PHE:N	2.46	0.49
1:C:312:GLU:HB2	1:C:315:GLU:HG3	1.94	0.49
1:D:315:GLU:O	1:D:319:VAL:HG23	2.12	0.49
1:E:9:PHE:CZ	1:E:55:LEU:HD11	2.47	0.49
1:G:58:ILE:O	1:G:58:ILE:HG13	2.11	0.49
1:F:168:VAL:HG21	1:F:189:LEU:HD13	1.94	0.49
1:H:219:ASP:OD2	1:H:222:LYS:HG2	2.11	0.49
1:C:130:TYR:CD1	1:C:289:ARG:HD3	2.47	0.49
1:H:31:LYS:HG2	1:H:123:TYR:CE1	2.48	0.49
1:E:271:ILE:HD12	1:E:271:ILE:H	1.77	0.48
1:F:172:GLY:O	1:F:177:GLY:HA3	2.13	0.48
1:A:192:VAL:HG22	1:A:211:ILE:HB	1.95	0.48
1:E:271:ILE:HG21	1:H:259:VAL:HG11	1.95	0.48
1:F:216:LEU:HD12	1:F:216:LEU:H	1.78	0.48
1:D:271:ILE:O	1:D:275:VAL:HG23	2.13	0.48
1:E:97:TYR:CZ	1:H:253:ARG:HB2	2.48	0.48
1:E:288:THR:HG22	1:E:291:ASP:H	1.79	0.48
1:H:138:LEU:HD11	1:H:300:ALA:HA	1.95	0.48
1:H:173:ILE:HD12	1:H:204:ALA:HB2	1.94	0.48
1:A:275:VAL:HG12	1:F:285:ILE:HB	1.95	0.48
1:A:288:THR:HG23	1:A:290:LYS:H	1.79	0.48
1:C:139:ASP:OD1	1:C:142:GLU:HG2	2.14	0.48
1:E:244:PHE:HE2	1:E:269:ILE:HG22	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:325:LYS:HB2	1:H:327:LYS:HG2	1.96	0.48
1:D:89:TYR:CZ	1:D:127:PRO:HD3	2.49	0.48
1:D:135:PRO:HD2	1:D:138:LEU:HD12	1.95	0.48
1:E:6:VAL:O	1:E:56:PRO:HA	2.14	0.48
1:D:309:GLU:OE1	1:D:322:ARG:NH2	2.37	0.48
1:G:288:THR:HG22	1:G:291:ASP:CG	2.33	0.47
1:D:172:GLY:O	1:D:177:GLY:HA3	2.13	0.47
1:E:288:THR:CG2	1:E:291:ASP:H	2.27	0.47
1:G:13:LEU:N	1:G:317:ASN:HD21	2.12	0.47
1:H:163:ARG:HA	1:H:163:ARG:HD3	1.72	0.47
1:A:152:THR:HG23	1:A:285:ILE:HA	1.97	0.47
1:D:178:HIS:CE1	1:D:179:ILE:HG13	2.49	0.47
1:F:288:THR:HG22	1:F:291:ASP:CG	2.34	0.47
1:G:82:ARG:HG2	1:G:134:ILE:CG2	2.44	0.47
1:G:70:VAL:HB	1:G:74:VAL:CG2	2.45	0.47
1:H:173:ILE:HA	1:H:177:GLY:HA3	1.97	0.47
1:C:38:CYS:SG	1:C:40:THR:OG1	2.70	0.47
1:E:92:CYS:SG	1:E:93:GLY:N	2.88	0.47
1:F:286:VAL:CG1	1:F:287:GLY:N	2.77	0.47
1:D:8:GLU:HG2	1:D:9:PHE:H	1.80	0.47
1:H:288:THR:CG2	1:H:291:ASP:H	2.25	0.47
1:A:244:PHE:HZ	1:A:269:ILE:HD11	1.81	0.46
1:A:315:GLU:O	1:A:319:VAL:HG23	2.15	0.46
1:A:10:LYS:HE2	1:A:10:LYS:HB3	1.70	0.46
1:G:173:ILE:O	1:G:173:ILE:HG13	2.14	0.46
1:G:230:GLY:HA3	1:G:252:LYS:HG3	1.97	0.46
1:H:20:ARG:HD3	1:H:123:TYR:CD2	2.50	0.46
1:H:70:VAL:HG13	1:H:79:VAL:HG23	1.98	0.46
1:H:67:VAL:HG12	1:H:81:ASP:O	2.16	0.46
1:E:311:ALA:O	1:E:334:LEU:HA	2.15	0.46
1:F:219:ASP:HB3	1:F:222:LYS:HB3	1.97	0.46
1:H:40:THR:HB	1:H:61:HIS:HE1	1.79	0.46
1:E:6:VAL:HG23	1:E:57:LEU:O	2.14	0.46
1:B:167:TRP:H	1:B:232:HIS:CE1	2.34	0.46
1:C:89:TYR:CZ	1:C:127:PRO:HD3	2.50	0.46
1:E:37:VAL:HG12	1:E:120:TYR:OH	2.15	0.46
1:G:61:HIS:HB3	1:G:112:GLY:O	2.16	0.46
1:B:199:GLU:OE1	1:B:199:GLU:N	2.42	0.45
1:B:89:TYR:CZ	1:B:127:PRO:HD3	2.52	0.45
1:C:297:ASP:O	1:C:301:ARG:HG3	2.15	0.45
1:D:316:ILE:HG23	1:D:317:ASN:ND2	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:58:ILE:N	1:H:117:ASP:OD1	2.36	0.45
1:A:316:ILE:HD13	1:A:334:LEU:HD11	1.99	0.45
1:G:136:ASP:OD1	1:G:136:ASP:N	2.49	0.45
1:H:175:GLY:O	1:H:178:HIS:ND1	2.48	0.45
1:G:87:TRP:CD1	1:G:286:VAL:HG21	2.52	0.45
1:F:38:CYS:HB3	1:F:331:ARG:CZ	2.47	0.45
1:G:181:LEU:HD21	1:G:209:ALA:HB2	1.98	0.45
1:B:42:LEU:HD12	1:B:323:MET:SD	2.56	0.45
1:B:97:TYR:CZ	1:D:253:ARG:HB2	2.51	0.45
1:F:181:LEU:HD13	1:F:191:VAL:HG11	1.98	0.45
1:H:89:TYR:CZ	1:H:127:PRO:HD3	2.52	0.45
1:H:175:GLY:O	1:H:178:HIS:CE1	2.69	0.45
1:D:36:GLY:H	1:D:62:GLU:HB2	1.82	0.45
1:H:171:TYR:HE2	1:H:246:GLN:HB3	1.81	0.45
1:F:31:LYS:O	1:F:31:LYS:HG3	2.16	0.45
1:A:202:LYS:HB2	1:A:202:LYS:HE2	1.83	0.45
1:G:140:PRO:HA	1:G:143:VAL:HG12	1.98	0.45
1:H:158:LYS:NZ	1:H:291:ASP:OD1	2.48	0.45
1:B:15:ILE:HD11	1:B:316:ILE:HG22	1.99	0.45
1:G:89:TYR:HB3	1:G:110:LEU:O	2.16	0.45
1:G:175:GLY:HA3	1:G:331:ARG:NE	2.29	0.45
1:H:286:VAL:HG22	1:H:287:GLY:H	1.82	0.45
1:H:321:GLU:O	1:H:325:LYS:HG2	2.17	0.45
1:B:222:LYS:HE2	1:B:226:ASP:OD2	2.17	0.44
1:C:37:VAL:HG13	1:C:120:TYR:OH	2.17	0.44
1:E:90:SER:HB3	1:E:127:PRO:HG2	1.99	0.44
1:D:127:PRO:O	1:D:131:VAL:HG13	2.18	0.44
1:E:246:GLN:O	1:E:250:SER:OG	2.34	0.44
1:E:319:VAL:HG11	1:E:332:ILE:HG21	2.00	0.44
1:F:53:PRO:HD3	1:F:114:TYR:CZ	2.52	0.44
1:F:173:ILE:HG12	1:F:195:ASP:HB2	1.99	0.44
1:G:2:LYS:HZ1	1:G:15:ILE:HG21	1.83	0.44
1:C:20:ARG:HH12	1:C:29:LEU:CD1	2.29	0.44
1:C:53:PRO:HD3	1:C:114:TYR:CZ	2.52	0.44
1:E:29:LEU:HD21	1:E:123:TYR:HB3	1.99	0.44
1:H:173:ILE:O	1:H:173:ILE:HG12	2.17	0.44
1:F:30:VAL:HG12	1:F:67:VAL:HG22	1.98	0.44
1:H:11:LYS:HA	1:H:11:LYS:HD2	1.61	0.44
1:H:31:LYS:HE3	1:H:123:TYR:CE1	2.53	0.44
1:D:57:LEU:HD11	1:D:113:GLY:O	2.17	0.44
1:E:1:MET:CE	1:E:121:ALA:HB1	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:315:GLU:O	1:E:319:VAL:HG22	2.18	0.44
1:G:245:GLU:O	1:G:249:GLN:HG3	2.18	0.44
1:H:245:GLU:O	1:H:249:GLN:HG3	2.17	0.44
1:B:149:ALA:HB3	1:B:179:ILE:HD12	1.99	0.44
1:E:40:THR:HB	1:E:61:HIS:HE1	1.83	0.44
1:B:28:VAL:HG12	1:B:70:VAL:HG12	2.00	0.44
1:E:313:LEU:HA	1:E:334:LEU:HD22	2.00	0.44
1:H:48:ASP:OD1	1:H:48:ASP:N	2.43	0.44
1:B:286:VAL:HG22	1:B:287:GLY:H	1.83	0.43
1:G:89:TYR:CE1	1:G:126:ALA:HA	2.53	0.43
1:G:296:LEU:HD23	1:G:296:LEU:HA	1.69	0.43
1:C:40:THR:OG1	1:C:61:HIS:NE2	2.45	0.43
1:C:178:HIS:CE1	1:C:179:ILE:HG13	2.53	0.43
1:G:82:ARG:CZ	1:G:140:PRO:HD3	2.48	0.43
1:B:181:LEU:HD23	1:B:182:GLN:H	1.84	0.43
1:C:137:ASN:OD1	1:C:137:ASN:N	2.52	0.43
1:E:170:ILE:HD12	1:E:181:LEU:HD23	1.99	0.43
1:G:147:LEU:HD22	1:G:292:MET:SD	2.59	0.43
1:G:151:VAL:HG11	1:G:292:MET:SD	2.58	0.43
1:G:328:ILE:HG12	1:G:332:ILE:HD11	1.99	0.43
1:H:320:PHE:C	1:H:324:GLU:OE2	2.57	0.43
1:A:322:ARG:HB3	1:A:328:ILE:HD11	2.00	0.43
1:G:112:GLY:HA2	1:G:116:VAL:CG1	2.47	0.43
1:A:75:LYS:O	1:A:78:LYS:NZ	2.45	0.43
1:A:87:TRP:O	1:A:111:ASN:HA	2.18	0.43
1:B:1:MET:SD	1:B:123:TYR:HB2	2.59	0.43
1:E:322:ARG:HG3	1:E:327:LYS:CE	2.47	0.43
1:B:167:TRP:CD2	1:B:228:VAL:HG21	2.53	0.43
1:E:42:LEU:HD23	1:E:42:LEU:HA	1.92	0.43
1:E:271:ILE:O	1:E:275:VAL:HG22	2.19	0.43
1:A:286:VAL:HG22	1:A:287:GLY:H	1.84	0.43
1:B:288:THR:HG22	1:B:290:LYS:N	2.29	0.43
1:B:36:GLY:HA2	1:B:333:VAL:HA	2.01	0.43
1:D:309:GLU:CD	1:D:322:ARG:HH22	2.18	0.43
1:F:220:PRO:O	1:F:224:ILE:HG13	2.19	0.43
1:C:231:VAL:O	1:C:252:LYS:HG2	2.19	0.43
1:F:149:ALA:HB2	1:F:331:ARG:NH2	2.34	0.43
1:G:40:THR:HG23	1:G:49:TRP:CH2	2.54	0.43
1:H:196:ILE:HB	1:H:216:LEU:HD21	1.99	0.43
1:H:256:THR:HG23	1:H:280:SER:HB2	2.01	0.43
1:A:137:ASN:OD1	1:A:137:ASN:N	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:152:THR:HG23	1:H:285:ILE:HA	2.00	0.42
1:G:286:VAL:HG12	1:G:287:GLY:N	2.34	0.42
1:F:196:ILE:HB	1:F:216:LEU:HD11	2.01	0.42
1:G:288:THR:HG23	1:G:290:LYS:H	1.85	0.42
1:H:124:CYS:SG	1:H:125:LYS:N	2.92	0.42
1:E:89:TYR:CZ	1:E:127:PRO:HD3	2.55	0.42
1:G:288:THR:CG2	1:G:291:ASP:H	2.32	0.42
1:G:328:ILE:CG2	1:G:332:ILE:HD12	2.50	0.42
1:H:49:TRP:CZ2	1:H:262:LEU:HD12	2.55	0.42
1:F:18:VAL:HG22	1:F:19:GLU:H	1.84	0.42
1:B:271:ILE:O	1:B:275:VAL:HG23	2.19	0.42
1:E:135:PRO:HD2	1:E:138:LEU:HD12	2.01	0.42
1:F:92:CYS:SG	1:F:93:GLY:N	2.92	0.42
1:F:139:ASP:OD2	1:F:141:VAL:HG12	2.20	0.42
1:G:246:GLN:O	1:G:250:SER:OG	2.32	0.42
1:H:200:LYS:O	1:H:203:LEU:HB3	2.20	0.42
1:B:183:TYR:HB3	1:B:187:MET:HE2	2.00	0.42
1:D:53:PRO:HD3	1:D:114:TYR:CZ	2.55	0.42
1:E:2:LYS:O	1:E:122:GLU:HG2	2.19	0.42
1:E:18:VAL:HG12	1:E:19:GLU:H	1.85	0.42
1:F:171:TYR:CE2	1:F:194:VAL:HG11	2.54	0.42
1:C:319:VAL:HG13	1:C:332:ILE:HD12	2.01	0.42
1:E:9:PHE:HZ	1:E:53:PRO:O	2.03	0.42
1:E:285:ILE:O	1:E:286:VAL:HG12	2.19	0.42
1:F:267:LEU:HD23	1:F:267:LEU:HA	1.91	0.42
1:G:40:THR:HB	1:G:61:HIS:CE1	2.49	0.42
1:G:221:VAL:HG23	1:G:246:GLN:HB3	2.02	0.42
1:H:4:ALA:HB2	1:H:313:LEU:HD21	2.02	0.42
1:B:88:LEU:HD21	1:B:103:GLU:HG3	2.00	0.42
1:C:23:LEU:HD21	1:C:28:VAL:N	2.35	0.42
1:E:244:PHE:CE2	1:E:269:ILE:HG22	2.55	0.42
1:G:328:ILE:HG23	1:G:332:ILE:HD12	2.01	0.42
1:C:183:TYR:O	1:C:187:MET:HG3	2.20	0.41
1:E:152:THR:HG23	1:E:285:ILE:HA	2.02	0.41
1:G:236:SER:HB2	1:G:257:LEU:HD11	2.02	0.41
1:G:309:GLU:CD	1:G:322:ARG:HH22	2.23	0.41
1:H:285:ILE:HD12	1:H:285:ILE:O	2.20	0.41
1:A:178:HIS:CE1	1:A:179:ILE:HG13	2.55	0.41
1:G:9:PHE:C	1:G:11:LYS:H	2.24	0.41
1:B:6:VAL:HG22	1:B:13:LEU:HG	2.01	0.41
1:D:160:SER:HB2	1:D:233:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:55:LEU:N	1:E:55:LEU:HD22	2.34	0.41
1:B:286:VAL:HG13	1:B:287:GLY:N	2.35	0.41
1:C:196:ILE:HB	1:C:216:LEU:HD23	2.03	0.41
1:E:72:LYS:HA	1:E:72:LYS:HD3	1.86	0.41
1:F:321:GLU:O	1:F:325:LYS:HG2	2.20	0.41
1:G:2:LYS:NZ	1:G:314:GLU:OE2	2.52	0.41
1:H:290:LYS:O	1:H:294:GLU:HG3	2.20	0.41
1:B:172:GLY:O	1:B:177:GLY:HA3	2.20	0.41
1:B:285:ILE:HB	1:D:275:VAL:HG12	2.01	0.41
1:G:248:TYR:CE2	1:G:270:PRO:HG2	2.55	0.41
1:A:40:THR:HB	1:A:61:HIS:CE1	2.50	0.41
1:E:173:ILE:O	1:E:173:ILE:HG22	2.21	0.41
1:F:245:GLU:O	1:F:249:GLN:HG3	2.20	0.41
1:G:3:ALA:HB2	1:G:121:ALA:HB2	2.03	0.41
1:H:29:LEU:HD21	1:H:123:TYR:HD2	1.85	0.41
1:A:38:CYS:HB3	1:A:331:ARG:NH2	2.36	0.41
1:E:37:VAL:HG22	1:E:332:ILE:O	2.21	0.41
1:G:84:GLY:HA3	1:G:147:LEU:HD11	2.03	0.41
1:C:139:ASP:HB2	1:C:140:PRO:HD2	2.03	0.41
1:C:282:LYS:HG2	1:G:280:SER:HB3	2.03	0.41
1:D:236:SER:HB2	1:D:257:LEU:HD11	2.03	0.41
1:E:170:ILE:HD13	1:E:180:ALA:HB1	2.02	0.41
1:E:286:VAL:HG13	1:E:287:GLY:N	2.36	0.41
1:F:325:LYS:HA	1:F:325:LYS:HD3	1.85	0.41
1:G:22:LYS:HB3	1:G:22:LYS:HE2	1.96	0.41
1:G:169:ALA:HB2	1:G:231:VAL:HG11	2.02	0.41
1:H:14:GLU:HB3	1:H:16:LYS:HG3	2.02	0.41
1:H:29:LEU:HD11	1:H:123:TYR:HB3	2.02	0.41
1:H:172:GLY:O	1:H:177:GLY:HA3	2.21	0.41
1:C:160:SER:HB2	1:C:233:ALA:HB1	2.02	0.40
1:G:29:LEU:HD23	1:G:29:LEU:HA	1.94	0.40
1:G:328:ILE:HG23	1:G:332:ILE:CD1	2.51	0.40
1:A:244:PHE:CG	1:A:267:LEU:HD13	2.56	0.40
1:A:319:VAL:HG13	1:A:332:ILE:HD12	2.02	0.40
1:B:275:VAL:CG1	1:D:285:ILE:HB	2.51	0.40
1:G:15:ILE:HG13	1:G:313:LEU:CD2	2.50	0.40
1:G:70:VAL:HB	1:G:74:VAL:HG21	2.03	0.40
1:H:236:SER:O	1:H:236:SER:OG	2.39	0.40
1:C:57:LEU:HD21	1:C:113:GLY:O	2.20	0.40
1:D:70:VAL:HG13	1:D:79:VAL:HG22	2.03	0.40
1:C:272:PHE:CD1	1:G:262:LEU:HB3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:273:ASP:O	1:E:277:ASN:ND2	2.51	0.40
1:H:322:ARG:NH2	1:H:327:LYS:O	2.52	0.40
1:B:27:GLU:OE2	1:B:125:LYS:HD2	2.21	0.40
1:D:87:TRP:O	1:D:111:ASN:HA	2.22	0.40
1:D:178:HIS:HA	1:D:181:LEU:HD12	2.04	0.40
1:E:195:ASP:OD1	1:E:196:ILE:HG12	2.22	0.40
1:F:27:GLU:OE2	1:F:125:LYS:HD2	2.21	0.40
1:G:137:ASN:ND2	1:G:300:ALA:O	2.55	0.40
1:H:58:ILE:HB	1:H:117:ASP:HB3	2.03	0.40
1:H:77:ILE:HD13	1:H:77:ILE:HA	1.93	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	334/339 (98%)	323 (97%)	10 (3%)	1 (0%)	41 61
1	B	334/339 (98%)	321 (96%)	12 (4%)	1 (0%)	41 61
1	C	334/339 (98%)	320 (96%)	12 (4%)	2 (1%)	25 46
1	D	334/339 (98%)	324 (97%)	9 (3%)	1 (0%)	41 61
1	E	334/339 (98%)	316 (95%)	15 (4%)	3 (1%)	17 34
1	F	334/339 (98%)	319 (96%)	14 (4%)	1 (0%)	41 61
1	G	334/339 (98%)	321 (96%)	12 (4%)	1 (0%)	41 61
1	H	334/339 (98%)	313 (94%)	20 (6%)	1 (0%)	41 61
All	All	2672/2712 (98%)	2557 (96%)	104 (4%)	11 (0%)	34 56

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	286	VAL
1	B	286	VAL
1	E	286	VAL
1	F	286	VAL
1	G	286	VAL
1	H	286	VAL
1	C	286	VAL
1	D	286	VAL
1	C	47	GLY
1	E	9	PHE
1	E	173	ILE

5.3.2 Protein sidechains [i](#)

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	265/268 (99%)	254 (96%)	11 (4%)	30 54
1	B	265/268 (99%)	257 (97%)	8 (3%)	41 65
1	C	265/268 (99%)	257 (97%)	8 (3%)	41 65
1	D	265/268 (99%)	261 (98%)	4 (2%)	65 82
1	E	265/268 (99%)	256 (97%)	9 (3%)	37 62
1	F	265/268 (99%)	261 (98%)	4 (2%)	65 82
1	G	265/268 (99%)	258 (97%)	7 (3%)	46 70
1	H	265/268 (99%)	255 (96%)	10 (4%)	33 59
All	All	2120/2144 (99%)	2059 (97%)	61 (3%)	42 67

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	10	LYS
1	A	11	LYS
1	A	42	LEU
1	A	54	LYS

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Mol	Chain	Res	Type
1	A	75	LYS
1	A	137	ASN
1	A	178	HIS
1	A	288	THR
1	A	305	ARG
1	A	322	ARG
1	B	8	GLU
1	B	11	LYS
1	B	54	LYS
1	B	72	LYS
1	B	137	ASN
1	B	163	ARG
1	B	197	SER
1	B	198	ASP
1	C	23	LEU
1	C	42	LEU
1	C	137	ASN
1	C	163	ARG
1	C	197	SER
1	C	202	LYS
1	C	273	ASP
1	C	305	ARG
1	D	78	LYS
1	D	232	HIS
1	D	286	VAL
1	D	322	ARG
1	E	1	MET
1	E	11	LYS
1	E	54	LYS
1	E	96	GLU
1	E	218	GLU
1	E	236	SER
1	E	241	LYS
1	E	327	LYS
1	E	331	ARG
1	F	19	GLU
1	F	92	CYS
1	F	320	PHE
1	F	321	GLU
1	G	29	LEU
1	G	82	ARG
1	G	136	ASP

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Mol	Chain	Res	Type
1	G	202	LYS
1	G	301	ARG
1	G	312	GLU
1	G	313	LEU
1	H	8	GLU
1	H	11	LYS
1	H	137	ASN
1	H	198	ASP
1	H	226	ASP
1	H	236	SER
1	H	242	LYS
1	H	305	ARG
1	H	327	LYS
1	H	331	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	227	GLN
1	C	293	GLN
1	F	111	ASN
1	G	7	ASN
1	G	293	GLN
1	H	7	ASN
1	H	39	HIS
1	H	317	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 16 ligands modelled in this entry, 16 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

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	336/339 (99%)	-0.16	0	100	100	26, 40, 60, 79	0
1	B	336/339 (99%)	-0.10	0	100	100	29, 47, 70, 91	0
1	C	336/339 (99%)	0.14	9 (2%)	54	49	29, 55, 82, 97	0
1	D	336/339 (99%)	0.10	6 (1%)	68	63	30, 51, 79, 99	0
1	E	336/339 (99%)	0.05	0	100	100	27, 54, 84, 100	0
1	F	336/339 (99%)	0.18	8 (2%)	59	54	32, 57, 84, 140	0
1	G	336/339 (99%)	0.77	34 (10%)	7	4	38, 75, 100, 118	0
1	H	336/339 (99%)	0.92	46 (13%)	3	2	46, 83, 113, 144	0
All	All	2688/2712 (99%)	0.24	103 (3%)	40	32	26, 56, 94, 144	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	14	GLU	6.5
1	G	316	ILE	6.1
1	H	316	ILE	5.6
1	H	3	ALA	5.5
1	G	313	LEU	5.1
1	H	172	GLY	4.5
1	H	123	TYR	4.5
1	H	13	LEU	4.4
1	H	79	VAL	4.2
1	G	29	LEU	4.1
1	H	334	LEU	3.9
1	H	15	ILE	3.8
1	H	332	ILE	3.8
1	C	316	ILE	3.8
1	G	3	ALA	3.7
1	H	40	THR	3.7

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Mol	Chain	Res	Type	RSRZ
1	F	73	GLY	3.6
1	H	12	ALA	3.6
1	H	173	ILE	3.6
1	H	1	MET	3.5
1	H	37	VAL	3.5
1	G	140	PRO	3.4
1	D	329	ASN	3.4
1	G	126	ALA	3.4
1	G	79	VAL	3.4
1	G	320	PHE	3.4
1	G	21	PRO	3.4
1	H	204	ALA	3.4
1	G	13	LEU	3.3
1	G	196	ILE	3.2
1	D	5	VAL	3.2
1	C	6	VAL	3.2
1	G	336	LEU	3.2
1	G	15	ILE	3.2
1	H	113	GLY	3.1
1	G	144	ALA	3.0
1	H	141	VAL	3.0
1	C	7	ASN	3.0
1	G	141	VAL	3.0
1	G	329	ASN	3.0
1	H	9	PHE	2.9
1	H	35	CYS	2.9
1	H	14	GLU	2.8
1	G	23	LEU	2.8
1	G	123	TYR	2.8
1	H	320	PHE	2.7
1	H	329	ASN	2.7
1	G	321	GLU	2.7
1	G	68	VAL	2.7
1	H	44	ALA	2.6
1	H	31	LYS	2.6
1	H	41	ASP	2.6
1	H	4	ALA	2.6
1	D	45	ALA	2.6
1	G	120	TYR	2.5
1	H	120	TYR	2.5
1	F	5	VAL	2.5
1	G	28	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	30	VAL	2.5
1	G	335	LYS	2.5
1	F	14	GLU	2.5
1	H	196	ILE	2.4
1	H	73	GLY	2.4
1	D	44	ALA	2.4
1	C	56	PRO	2.4
1	D	3	ALA	2.4
1	C	13	LEU	2.4
1	F	120	TYR	2.4
1	H	140	PRO	2.3
1	H	2	LYS	2.3
1	H	112	GLY	2.3
1	D	73	GLY	2.3
1	H	237	VAL	2.3
1	H	122	GLU	2.3
1	F	45	ALA	2.3
1	G	308	VAL	2.3
1	H	321	GLU	2.3
1	F	4	ALA	2.3
1	H	32	ILE	2.3
1	G	70	VAL	2.2
1	G	31	LYS	2.2
1	H	57	LEU	2.2
1	H	176	LEU	2.2
1	C	45	ALA	2.2
1	G	4	ALA	2.2
1	G	333	VAL	2.2
1	G	116	VAL	2.2
1	H	175	GLY	2.2
1	H	20	ARG	2.2
1	C	15	ILE	2.2
1	C	329	ASN	2.2
1	H	83	VAL	2.2
1	H	313	LEU	2.1
1	G	76	SER	2.1
1	G	66	ILE	2.1
1	G	85	ILE	2.1
1	F	35	CYS	2.1
1	H	5	VAL	2.1
1	C	116	VAL	2.1
1	F	58	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	G	1	MET	2.0
1	H	335	LYS	2.0
1	H	119	GLY	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 monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	H	401	1/1	0.72	0.17	145,145,145,145	0
2	ZN	F	401	1/1	0.94	0.12	71,71,71,71	0
2	ZN	E	401	1/1	0.94	0.10	62,62,62,62	0
2	ZN	G	401	1/1	0.95	0.11	95,95,95,95	0
2	ZN	A	401	1/1	0.98	0.10	50,50,50,50	0
2	ZN	B	401	1/1	0.98	0.14	57,57,57,57	0
2	ZN	G	400	1/1	0.98	0.14	46,46,46,46	0
2	ZN	C	400	1/1	0.98	0.09	53,53,53,53	0
2	ZN	H	400	1/1	0.98	0.12	59,59,59,59	0
2	ZN	C	401	1/1	0.98	0.12	70,70,70,70	0
2	ZN	D	400	1/1	0.99	0.11	44,44,44,44	0
2	ZN	D	401	1/1	0.99	0.06	55,55,55,55	0
2	ZN	E	400	1/1	0.99	0.10	57,57,57,57	0
2	ZN	A	400	1/1	0.99	0.09	32,32,32,32	0
2	ZN	F	400	1/1	0.99	0.13	41,41,41,41	0
2	ZN	B	400	1/1	1.00	0.14	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.