



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2022 – 04:07 PM JST

PDB ID : 7W08
Title : Itaconate inducible LysR-Type Transcriptional regulator (ITCR) in APO form, Space group P1.
Authors : Sun, P.K.; Wang, B.; Li, X.J.
Deposited on : 2021-11-17
Resolution : 3.25 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	:	2.31.2
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)	:	2.31.2

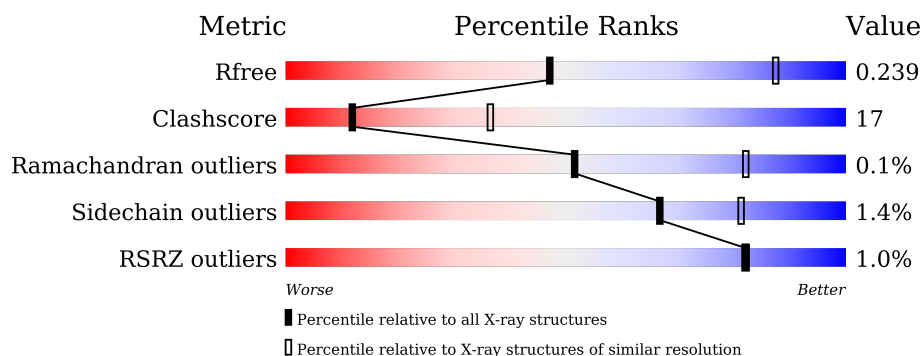
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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	1191 (3.30-3.22)
Clashscore	141614	1251 (3.30-3.22)
Ramachandran outliers	138981	1229 (3.30-3.22)
Sidechain outliers	138945	1228 (3.30-3.22)
RSRZ outliers	127900	1154 (3.30-3.22)

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	292	<div> <div>%</div> <div> <div></div> <div>47%</div> <div>23%</div> <div>30%</div> </div> </div>
1	B	292	<div> <div></div> <div>48%</div> <div>21%</div> <div>•</div> <div>31%</div> </div>
1	C	292	<div> <div>%</div> <div> <div></div> <div>46%</div> <div>24%</div> <div>•</div> <div>29%</div> </div> </div>
1	D	292	<div> <div></div> <div>48%</div> <div>21%</div> <div>•</div> <div>30%</div> </div>
1	E	292	<div> <div>%</div> <div> <div></div> <div>50%</div> <div>20%</div> <div>•</div> <div>29%</div> </div> </div>
1	F	292	<div> <div>%</div> <div> <div></div> <div>44%</div> <div>25%</div> <div>•</div> <div>30%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	292	<div><div><div>%</div><div><div></div><div>49%</div><div>21%</div><div>29%</div></div><div></div></div></div>
1	H	292	<div><div><div>%</div><div><div></div><div>49%</div><div>20%</div><div>29%</div></div><div></div></div></div>

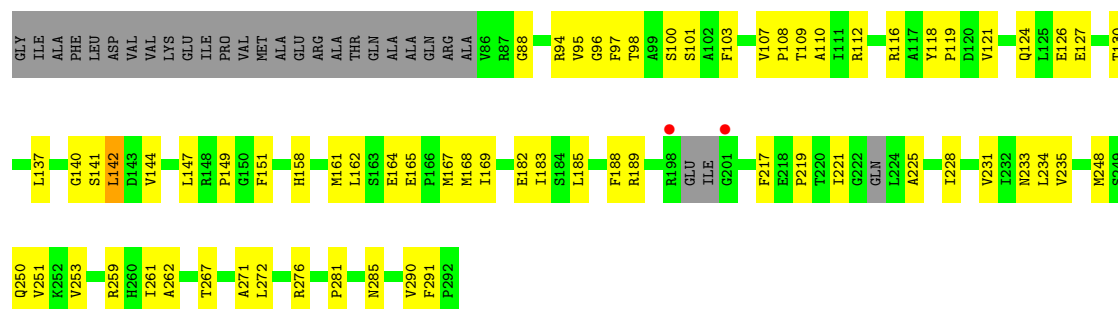
2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12198 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.

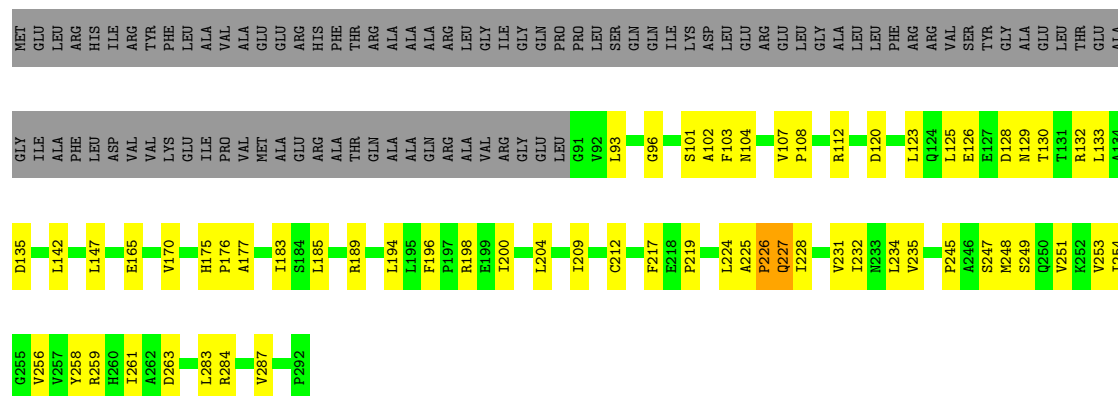
- Molecule 1 is a protein called Transcriptional regulator, LysR family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	207	Total	C	N	O	S	0	0	0
			1525	972	262	284	7			
1	D	204	Total	C	N	O	S	0	0	0
			1521	971	258	285	7			
1	A	204	Total	C	N	O	S	0	0	0
			1469	939	247	276	7			
1	B	202	Total	C	N	O	S	0	0	0
			1512	967	257	281	7			
1	E	207	Total	C	N	O	S	0	0	0
			1518	971	259	281	7			
1	F	205	Total	C	N	O	S	0	1	0
			1561	996	264	294	7			
1	G	207	Total	C	N	O	S	0	0	0
			1544	983	264	290	7			
1	H	207	Total	C	N	O	S	0	0	0
			1548	989	270	282	7			



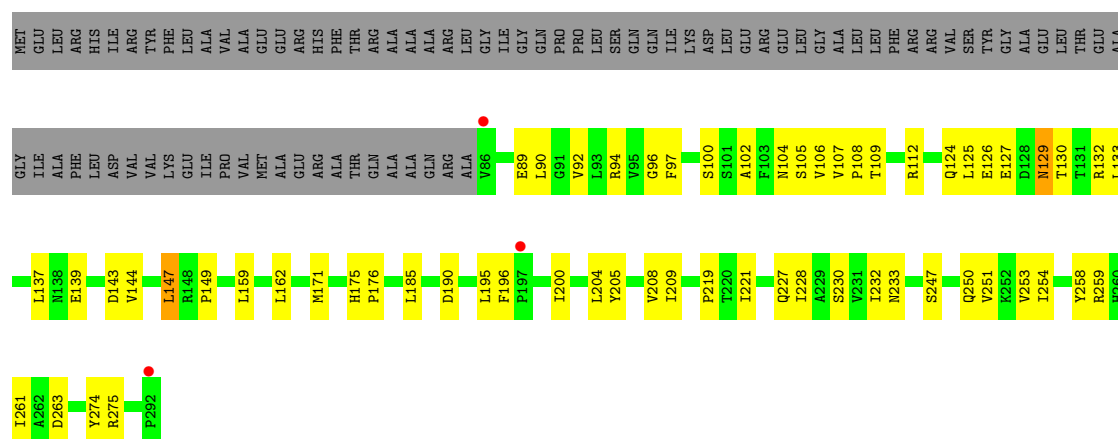
- Molecule 1: Transcriptional regulator, LysR family

Chain B: 48% 21% 31%



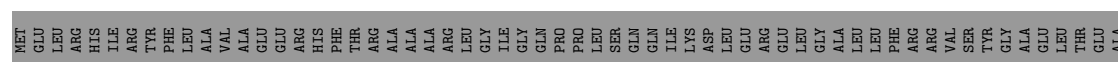
- Molecule 1: Transcriptional regulator, LysR family

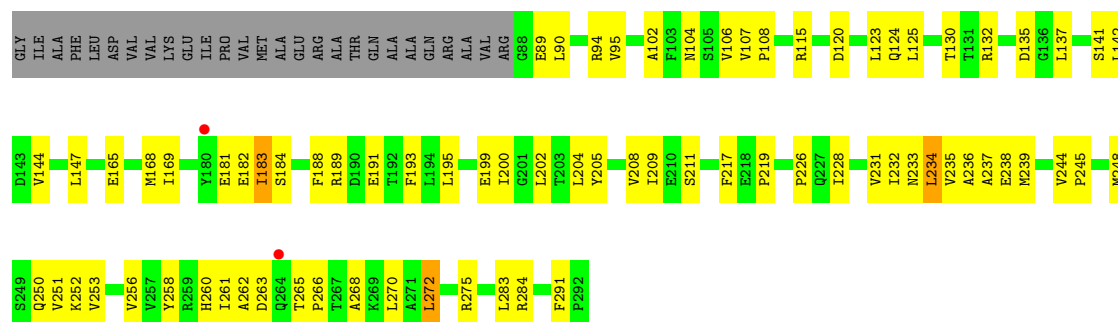
Chain E: 50% 20% 29%



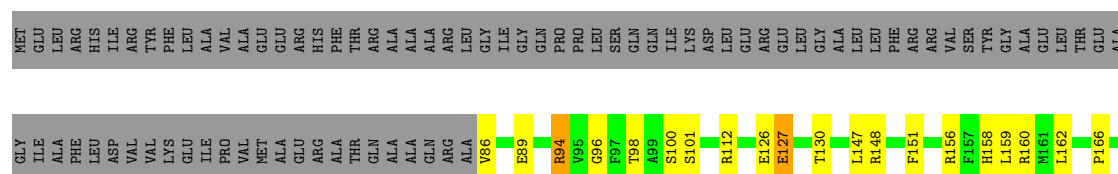
- Molecule 1: Transcriptional regulator, LysR family

Chain F: 44% 25% 30%

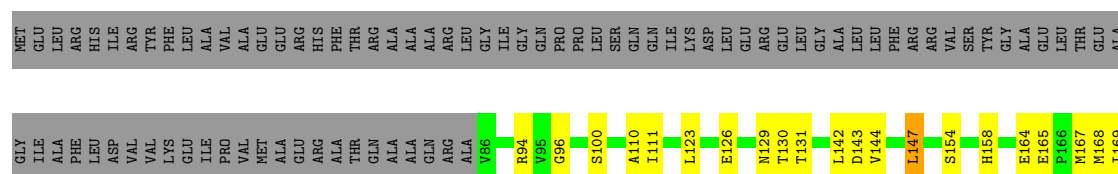




• Molecule 1: Transcriptional regulator, LysR family



• Molecule 1: Transcriptional regulator, LysR family



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.59Å 73.52Å 104.03Å 94.24° 105.67° 111.13°	Depositor
Resolution (Å)	32.96 – 3.25 40.37 – 3.25	Depositor EDS
% Data completeness (in resolution range)	91.4 (32.96-3.25) 87.2 (40.37-3.25)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.18_3855	Depositor
R, R_{free}	0.199 , 0.239 0.199 , 0.239	Depositor DCC
R_{free} test set	2000 reflections (7.75%)	wwPDB-VP
Wilson B-factor (Å ²)	51.1	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12198	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% 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

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	0.59	0/1495	0.87	3/2038 (0.1%)
1	B	0.63	0/1542	0.96	2/2101 (0.1%)
1	C	0.65	1/1554 (0.1%)	0.89	1/2119 (0.0%)
1	D	0.67	0/1551	0.97	3/2114 (0.1%)
1	E	0.57	0/1549	0.86	1/2114 (0.0%)
1	F	0.71	0/1592	0.96	5/2167 (0.2%)
1	G	0.72	2/1575 (0.1%)	0.90	4/2145 (0.2%)
1	H	0.72	0/1578	0.99	8/2149 (0.4%)
All	All	0.66	3/12436 (0.0%)	0.93	27/16947 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	245	PRO	N-CA	12.09	1.67	1.47
1	C	212	CYS	CB-SG	-5.47	1.72	1.81
1	G	244	VAL	C-N	5.20	1.44	1.34

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	226	PRO	N-CA-CB	12.59	118.41	103.30
1	D	283	LEU	CB-CG-CD1	-10.75	92.72	111.00
1	H	225	ALA	CB-CA-C	-7.90	98.25	110.10
1	D	168	MET	CB-CG-SD	-7.60	89.60	112.40
1	H	223	GLN	CB-CA-C	-7.41	95.58	110.40
1	H	224	LEU	N-CA-C	-7.10	91.83	111.00
1	C	239	MET	CB-CG-SD	-7.06	91.23	112.40
1	H	225	ALA	N-CA-CB	6.70	119.47	110.10
1	G	162	LEU	CB-CG-CD2	-6.51	99.94	111.00
1	H	225	ALA	N-CA-C	-6.48	93.50	111.00
1	A	168	MET	CG-SD-CE	6.43	110.49	100.20
1	H	202	LEU	CB-CG-CD2	-6.24	100.39	111.00
1	G	245	PRO	CA-N-CD	-6.11	102.94	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	LEU	CA-CB-CG	6.10	129.33	115.30
1	F	183	ILE	CG1-CB-CG2	-6.05	98.09	111.40
1	G	245	PRO	N-CA-C	-5.95	96.63	112.10
1	E	147	LEU	CB-CG-CD2	-5.93	100.92	111.00
1	H	198	ARG	NE-CZ-NH2	5.68	123.14	120.30
1	F	272	LEU	CB-CG-CD1	-5.66	101.39	111.00
1	D	226	PRO	N-CA-CB	5.63	110.06	103.30
1	B	227	GLN	CB-CA-C	-5.50	99.40	110.40
1	H	224	LEU	CB-CA-C	-5.50	99.75	110.20
1	F	234	LEU	CB-CG-CD1	-5.47	101.70	111.00
1	A	167	MET	CG-SD-CE	5.38	108.81	100.20
1	F	90	LEU	CB-CG-CD2	-5.28	102.02	111.00
1	G	94	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	F	252	LYS	CD-CE-NZ	5.10	123.43	111.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1469	0	1391	50	0
1	B	1512	0	1495	63	0
1	C	1525	0	1493	69	0
1	D	1521	0	1491	56	0
1	E	1518	0	1468	49	0
1	F	1561	0	1550	69	0
1	G	1544	0	1503	48	0
1	H	1548	0	1546	68	0
All	All	12198	0	11937	417	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 17.

All (417) 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:G:245:PRO:N	1:G:245:PRO:CA	1.67	1.47
1:D:225:ALA:HB1	1:B:225:ALA:HB1	1.24	1.16
1:H:225:ALA:CB	1:H:226:PRO:HD2	1.70	1.16
1:H:181:GLU:HA	1:H:259:ARG:HH22	1.17	1.05
1:H:225:ALA:HB1	1:H:226:PRO:HD2	1.38	1.04
1:F:199:GLU:OE2	1:F:231:VAL:HG12	1.59	1.00
1:C:139:GLU:OE1	1:B:198:ARG:HG3	1.61	0.99
1:H:225:ALA:CB	1:H:226:PRO:CD	2.40	0.99
1:C:148:ARG:HD2	1:C:204:LEU:CD1	1.94	0.97
1:H:225:ALA:HB3	1:H:226:PRO:HD2	1.43	0.97
1:D:202:LEU:HD21	1:A:140:GLY:HA3	1.46	0.96
1:C:124:GLN:HE22	1:B:225:ALA:HB2	1.29	0.96
1:H:225:ALA:HB1	1:H:226:PRO:CD	1.95	0.95
1:F:130:THR:HG23	1:F:147:LEU:HB2	1.49	0.94
1:D:225:ALA:HB1	1:B:225:ALA:CB	1.97	0.93
1:C:124:GLN:NE2	1:B:225:ALA:HB2	1.85	0.90
1:H:130:THR:HG23	1:H:147:LEU:HB2	1.55	0.88
1:E:149:PRO:HB3	1:E:159:LEU:HD21	1.54	0.88
1:C:92:VAL:HG21	1:B:224:LEU:CD1	2.04	0.87
1:G:253:VAL:HA	1:H:250:GLN:HG2	1.57	0.87
1:E:130:THR:HG23	1:E:147:LEU:HB2	1.57	0.86
1:D:227:GLN:HE21	1:A:141:SER:HB2	1.41	0.86
1:F:228:ILE:HD11	1:F:234:LEU:HD11	1.56	0.85
1:A:164:GLU:OE1	1:A:267:THR:OG1	1.93	0.83
1:A:130:THR:HG23	1:A:147:LEU:HB2	1.60	0.83
1:H:181:GLU:HA	1:H:259:ARG:NH2	1.94	0.82
1:D:225:ALA:CB	1:B:225:ALA:HB1	2.09	0.80
1:E:124:GLN:NE2	1:F:226:PRO:HG3	1.98	0.79
1:B:200:ILE:HG23	1:B:204:LEU:HD23	1.64	0.79
1:E:129:ASN:O	1:E:133:LEU:HG	1.83	0.78
1:B:132:ARG:HA	1:B:135:ASP:HB2	1.63	0.78
1:D:225:ALA:CB	1:B:225:ALA:CB	2.62	0.77
1:E:92:VAL:HG21	1:H:224:LEU:HD13	1.65	0.77
1:C:92:VAL:HG21	1:B:224:LEU:HD11	1.66	0.76
1:C:139:GLU:HG2	1:C:141:SER:HB3	1.65	0.76
1:D:148:ARG:HD2	1:D:204:LEU:HD13	1.69	0.75
1:B:196:PHE:CD2	1:B:231:VAL:HG22	2.22	0.75
1:A:272:LEU:HD23	1:A:291:PHE:HE1	1.51	0.75
1:E:124:GLN:HE22	1:F:226:PRO:HG3	1.50	0.75
1:G:245:PRO:N	1:G:245:PRO:C	2.40	0.74
1:D:227:GLN:NE2	1:A:141:SER:HB2	2.02	0.74
1:G:98:THR:HG1	1:G:101:SER:HG	0.74	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:183:ILE:HG22	1:F:184:SER:H	1.52	0.73
1:G:196:PHE:HE2	1:G:205:TYR:HB2	1.54	0.72
1:E:261:ILE:HG22	1:E:263:ASP:H	1.55	0.72
1:F:209:ILE:HD13	1:F:219:PRO:HG3	1.72	0.72
1:B:198:ARG:CZ	1:B:227:GLN:HB3	2.19	0.72
1:B:120:ASP:OD2	1:H:282:VAL:HG23	1.91	0.71
1:G:100:SER:HB3	1:G:228:ILE:HD12	1.72	0.70
1:G:196:PHE:CE2	1:G:205:TYR:HB2	2.26	0.70
1:E:137:LEU:O	1:E:275:ARG:NH1	2.25	0.70
1:H:129:ASN:ND2	1:H:131:THR:OG1	2.24	0.70
1:A:169:ILE:HB	1:A:261:ILE:HD11	1.74	0.69
1:C:100:SER:HB3	1:C:228:ILE:HG13	1.75	0.69
1:D:94:ARG:NH1	1:E:190:ASP:OD1	2.26	0.69
1:G:171:MET:HE1	1:G:177:ALA:HB3	1.75	0.68
1:F:199:GLU:OE2	1:F:231:VAL:CG1	2.37	0.68
1:G:212:CYS:SG	1:G:219:PRO:HG3	2.34	0.67
1:E:149:PRO:HB3	1:E:159:LEU:CD2	2.24	0.66
1:C:272:LEU:HD23	1:C:291:PHE:HE2	1.61	0.66
1:C:148:ARG:CD	1:C:204:LEU:HD11	2.26	0.66
1:C:148:ARG:HD2	1:C:204:LEU:HD11	1.77	0.65
1:C:153:GLY:O	1:C:156:ARG:HD3	1.97	0.65
1:C:265:THR:N	1:C:266:PRO:HD3	2.11	0.65
1:D:147:LEU:HD13	1:D:151:PHE:HE2	1.61	0.64
1:E:112:ARG:HD3	1:F:238:GLU:HG3	1.79	0.64
1:B:212:CYS:HB3	1:B:219:PRO:HD3	1.80	0.64
1:C:92:VAL:HG21	1:B:224:LEU:HD13	1.79	0.64
1:C:148:ARG:CD	1:C:204:LEU:CD1	2.74	0.63
1:C:185:LEU:HG	1:C:261:ILE:HG21	1.80	0.63
1:G:265:THR:O	1:G:265:THR:OG1	2.08	0.63
1:H:225:ALA:O	1:H:226:PRO:C	2.36	0.63
1:B:130:THR:HG23	1:B:147:LEU:HB2	1.79	0.63
1:C:124:GLN:HE22	1:B:225:ALA:CB	2.06	0.63
1:B:103:PHE:O	1:B:247:SER:HB3	1.99	0.63
1:E:139:GLU:OE2	1:H:198:ARG:NH2	2.23	0.63
1:B:198:ARG:NH1	1:B:227:GLN:HB3	2.14	0.63
1:F:204:LEU:HD21	1:F:245:PRO:HG3	1.80	0.62
1:F:245:PRO:HD2	1:F:248:MET:HG3	1.81	0.62
1:E:102:ALA:O	1:F:233:ASN:ND2	2.32	0.62
1:F:200:ILE:HG22	1:F:204:LEU:HB3	1.82	0.62
1:E:196:PHE:HZ	1:E:204:LEU:HB3	1.64	0.61
1:F:189:ARG:HA	1:F:217:PHE:CE2	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:132:ARG:HG2	1:F:132:ARG:HH21	1.65	0.60
1:F:94:ARG:HG3	1:F:124:GLN:HB3	1.82	0.60
1:H:164:GLU:HG2	1:H:267:THR:OG1	2.01	0.60
1:H:199:GLU:HG3	1:H:229:ALA:HB1	1.84	0.60
1:C:148:ARG:HD2	1:C:204:LEU:HD13	1.80	0.60
1:G:160:ARG:HH11	1:G:160:ARG:HG3	1.67	0.60
1:H:200:ILE:HG22	1:H:204:LEU:HB3	1.83	0.60
1:C:92:VAL:CG2	1:B:224:LEU:CD1	2.79	0.59
1:H:94:ARG:NH2	1:H:143:ASP:OD1	2.35	0.59
1:H:96:GLY:HA2	1:H:126:GLU:O	2.01	0.59
1:D:258:TYR:O	1:D:259:ARG:HD2	2.03	0.59
1:C:144:VAL:CG2	1:C:272:LEU:HD11	2.33	0.59
1:H:158:HIS:CE1	1:H:276:ARG:HA	2.38	0.59
1:E:149:PRO:CB	1:E:159:LEU:HD21	2.29	0.59
1:D:224:LEU:CB	1:A:124:GLN:HG3	2.33	0.59
1:B:224:LEU:HD23	1:B:224:LEU:H	1.68	0.58
1:H:100:SER:HB3	1:H:199:GLU:O	2.03	0.58
1:A:96:GLY:HA2	1:A:126:GLU:O	2.02	0.58
1:H:181:GLU:CA	1:H:259:ARG:HH22	2.05	0.58
1:D:158:HIS:CE1	1:D:276:ARG:HH21	2.21	0.58
1:H:251:VAL:HG12	1:H:253:VAL:HG23	1.85	0.58
1:A:219:PRO:O	1:A:221:ILE:HG13	2.04	0.58
1:C:95:VAL:HB	1:C:125:LEU:HD23	1.85	0.58
1:E:129:ASN:ND2	1:E:200:ILE:O	2.37	0.58
1:D:118:TYR:OH	1:D:285:ASN:ND2	2.37	0.57
1:E:139:GLU:CD	1:H:198:ARG:HH21	2.05	0.57
1:D:247:SER:O	1:D:250:GLN:NE2	2.30	0.57
1:D:107:VAL:HB	1:D:108:PRO:HD3	1.86	0.57
1:A:109:THR:HG23	1:A:112:ARG:HH21	1.70	0.57
1:D:202:LEU:HD21	1:A:140:GLY:CA	2.27	0.57
1:B:133:LEU:HD22	1:B:142:LEU:HD11	1.86	0.57
1:H:276:ARG:HG2	1:H:276:ARG:HH11	1.69	0.57
1:A:116:ARG:O	1:A:119:PRO:HD3	2.05	0.56
1:D:175:HIS:ND1	1:D:176:PRO:HD2	2.21	0.56
1:G:171:MET:HE3	1:G:175:HIS:HB3	1.87	0.56
1:G:208:VAL:HG13	1:G:243:ILE:HD12	1.87	0.56
1:G:249:SER:O	1:G:249:SER:OG	2.21	0.56
1:B:227:GLN:O	1:B:227:GLN:HG3	2.06	0.56
1:G:130:THR:HG23	1:G:147:LEU:HB2	1.86	0.56
1:G:197:PRO:HA	1:G:224:LEU:HD13	1.88	0.56
1:G:285:ASN:O	1:G:289:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:212:CYS:HB2	1:H:219:PRO:HD3	1.88	0.56
1:C:165:GLU:HG3	1:C:268:ALA:HB3	1.88	0.56
1:D:164:GLU:OE1	1:D:269:LYS:HE3	2.06	0.56
1:B:251:VAL:HG12	1:B:253:VAL:HG23	1.86	0.56
1:F:183:ILE:HG22	1:F:184:SER:N	2.21	0.56
1:A:225:ALA:HB3	1:A:231:VAL:HG23	1.87	0.55
1:H:228:ILE:HD12	1:H:234:LEU:HD11	1.88	0.55
1:C:92:VAL:CG2	1:B:224:LEU:HD11	2.34	0.55
1:E:100:SER:HB3	1:E:228:ILE:HD12	1.87	0.55
1:E:233:ASN:HB3	1:F:102:ALA:HB1	1.88	0.55
1:H:158:HIS:O	1:H:273:ALA:HA	2.06	0.55
1:C:197:PRO:CB	1:C:224:LEU:HD11	2.36	0.55
1:D:280:SER:HB3	1:D:283:LEU:HB3	1.87	0.55
1:A:100:SER:HB3	1:A:228:ILE:HD12	1.87	0.55
1:E:232:ILE:HD13	1:E:258:TYR:OH	2.07	0.55
1:C:118:TYR:CE2	1:C:285:ASN:HB3	2.42	0.55
1:C:144:VAL:HG22	1:C:272:LEU:HD11	1.89	0.55
1:F:200:ILE:CG2	1:F:204:LEU:HB3	2.36	0.55
1:B:129:ASN:O	1:B:133:LEU:HG	2.06	0.54
1:G:183:ILE:HG21	1:G:188:PHE:HE1	1.71	0.54
1:B:185:LEU:HG	1:B:261:ILE:HG21	1.89	0.54
1:E:129:ASN:HB2	1:E:132:ARG:H	1.72	0.54
1:C:139:GLU:OE1	1:B:198:ARG:CG	2.47	0.54
1:C:231:VAL:O	1:C:235:VAL:HG23	2.08	0.54
1:H:180:TYR:O	1:H:259:ARG:NH1	2.30	0.54
1:B:249:SER:HA	1:B:258:TYR:CE2	2.43	0.54
1:E:97:PHE:CE1	1:E:127:GLU:HB2	2.42	0.54
1:E:107:VAL:HB	1:E:108:PRO:HD3	1.90	0.54
1:E:247:SER:O	1:E:250:GLN:HG3	2.08	0.54
1:A:158:HIS:CE1	1:A:276:ARG:HA	2.43	0.54
1:A:250:GLN:OE1	1:B:254:ILE:HG13	2.07	0.53
1:H:144:VAL:HG22	1:H:272:LEU:HD11	1.90	0.53
1:B:175:HIS:CD2	1:B:176:PRO:HD2	2.44	0.53
1:G:175:HIS:CD2	1:G:176:PRO:HD2	2.43	0.53
1:H:167:MET:HG3	1:H:245:PRO:HA	1.90	0.53
1:C:265:THR:O	1:C:265:THR:OG1	2.22	0.53
1:E:195:LEU:HD21	1:E:208:VAL:HG11	1.91	0.53
1:A:118:TYR:HB3	1:A:121:VAL:HG23	1.90	0.53
1:G:158:HIS:O	1:G:273:ALA:HA	2.09	0.53
1:D:92:VAL:HG23	1:D:122:ARG:O	2.08	0.53
1:B:198:ARG:HD2	1:B:198:ARG:N	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:253:VAL:O	1:D:256:VAL:HG22	2.09	0.52
1:F:137:LEU:O	1:F:275:ARG:NH1	2.42	0.52
1:H:189:ARG:HB2	1:H:217:PHE:CG	2.44	0.52
1:D:132:ARG:HG2	1:D:132:ARG:HH21	1.74	0.52
1:A:107:VAL:HB	1:A:108:PRO:HD3	1.92	0.52
1:C:130:THR:HG23	1:C:147:LEU:HB2	1.91	0.52
1:A:149:PRO:HD3	1:A:271:ALA:HB3	1.91	0.52
1:D:128:ASP:OD1	1:D:129:ASN:N	2.42	0.52
1:C:112:ARG:HG3	1:D:238:GLU:HB2	1.90	0.52
1:A:189:ARG:HA	1:A:217:PHE:CE2	2.45	0.52
1:B:235:VAL:HG12	1:B:256:VAL:HG12	1.92	0.52
1:G:217:PHE:HE2	1:G:219:PRO:HB3	1.74	0.52
1:D:169:ILE:HG22	1:D:259:ARG:O	2.09	0.52
1:A:183:ILE:HG23	1:A:188:PHE:HE2	1.75	0.52
1:D:225:ALA:CB	1:B:225:ALA:HB3	2.39	0.52
1:B:198:ARG:N	1:B:198:ARG:CD	2.72	0.52
1:B:198:ARG:NE	1:B:227:GLN:HB3	2.24	0.52
1:C:157:PHE:HB3	1:C:274:TYR:O	2.09	0.51
1:C:160:ARG:HG3	1:C:160:ARG:HH11	1.74	0.51
1:A:127:GLU:OE1	1:B:228:ILE:HG23	2.10	0.51
1:A:149:PRO:HD3	1:A:271:ALA:CB	2.40	0.51
1:H:158:HIS:ND1	1:H:276:ARG:HA	2.24	0.51
1:H:250:GLN:O	1:H:252:LYS:N	2.43	0.51
1:D:104:ASN:OD1	1:D:106:VAL:HG13	2.10	0.51
1:A:88:GLY:HA3	1:A:281:PRO:HB2	1.93	0.51
1:E:112:ARG:HD3	1:F:238:GLU:CG	2.41	0.51
1:F:195:LEU:HD21	1:F:208:VAL:CG1	2.41	0.51
1:A:189:ARG:HA	1:A:217:PHE:CD2	2.45	0.51
1:F:132:ARG:HA	1:F:135:ASP:HB3	1.92	0.51
1:F:193:PHE:HB3	1:F:195:LEU:HD13	1.93	0.51
1:B:251:VAL:O	1:B:253:VAL:N	2.41	0.51
1:F:168:MET:HE3	1:F:260:HIS:NE2	2.26	0.51
1:C:125:LEU:HD11	1:D:239:MET:SD	2.50	0.51
1:D:228:ILE:HD11	1:D:234:LEU:HD11	1.93	0.51
1:D:244:VAL:HB	1:D:248:MET:HG3	1.92	0.50
1:B:177:ALA:HB1	1:B:183:ILE:HD11	1.93	0.50
1:H:225:ALA:HB3	1:H:226:PRO:CD	2.24	0.50
1:C:167:MET:HG3	1:C:266:PRO:O	2.12	0.50
1:D:208:VAL:HG13	1:D:243:ILE:HD12	1.94	0.50
1:A:259:ARG:HG3	1:A:259:ARG:HH11	1.76	0.50
1:C:113:ALA:O	1:C:116:ARG:HB3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:92:VAL:HG21	1:H:224:LEU:CD1	2.38	0.50
1:G:160:ARG:HG3	1:G:160:ARG:NH1	2.27	0.50
1:B:96:GLY:HA2	1:B:126:GLU:O	2.12	0.50
1:F:94:ARG:CG	1:F:124:GLN:HB3	2.41	0.50
1:H:213:ARG:HE	1:H:218:GLU:CD	2.14	0.49
1:D:147:LEU:HD13	1:D:151:PHE:CE2	2.46	0.49
1:C:151:PHE:CD1	1:C:151:PHE:C	2.86	0.49
1:E:205:TYR:O	1:E:209:ILE:HG12	2.11	0.49
1:F:107:VAL:HB	1:F:108:PRO:HD3	1.94	0.49
1:A:185:LEU:HD11	1:A:261:ILE:HD13	1.94	0.49
1:F:165:GLU:HG3	1:F:268:ALA:HB3	1.93	0.49
1:F:183:ILE:O	1:F:262:ALA:N	2.27	0.49
1:H:165:GLU:OE1	1:H:247:SER:OG	2.23	0.49
1:F:232:ILE:HD13	1:F:258:TYR:OH	2.12	0.49
1:C:236:ALA:O	1:D:112:ARG:NH2	2.46	0.49
1:G:170:VAL:HG13	1:G:235:VAL:HG21	1.94	0.49
1:F:94:ARG:HD2	1:F:141:SER:O	2.12	0.49
1:F:211:SER:OG	1:F:266:PRO:HD3	2.12	0.49
1:C:103:PHE:HA	1:D:233:ASN:HD21	1.77	0.49
1:C:115:ARG:HD2	1:C:123:LEU:HD12	1.93	0.49
1:H:247:SER:O	1:H:250:GLN:OE1	2.31	0.49
1:C:89:GLU:HG2	1:C:118:TYR:CE1	2.47	0.49
1:C:209:ILE:HD13	1:C:209:ILE:N	2.27	0.49
1:D:166:PRO:HA	1:D:267:THR:HG22	1.94	0.49
1:B:232:ILE:HD13	1:B:258:TYR:OH	2.13	0.49
1:H:189:ARG:HA	1:H:217:PHE:CE2	2.48	0.49
1:G:209:ILE:O	1:G:213:ARG:HG3	2.13	0.48
1:E:254:ILE:HG13	1:F:250:GLN:OE1	2.13	0.48
1:F:265:THR:HB	1:F:266:PRO:CD	2.43	0.48
1:F:251:VAL:HG12	1:F:253:VAL:HG23	1.95	0.48
1:B:189:ARG:HA	1:B:217:PHE:CD2	2.49	0.48
1:G:127:GLU:CD	1:H:228:ILE:HG23	2.34	0.48
1:G:170:VAL:HG22	1:G:235:VAL:HG11	1.95	0.48
1:C:252:LYS:O	1:D:250:GLN:HB3	2.13	0.48
1:A:231:VAL:O	1:A:235:VAL:HG23	2.14	0.48
1:H:276:ARG:HG2	1:H:276:ARG:NH1	2.28	0.47
1:C:183:ILE:HD13	1:C:183:ILE:HA	1.71	0.47
1:E:105:SER:O	1:E:108:PRO:HD2	2.14	0.47
1:G:175:HIS:CG	1:G:176:PRO:HD2	2.49	0.47
1:H:202:LEU:HA	1:H:202:LEU:HD23	1.47	0.47
1:C:188:PHE:CE2	1:C:241:VAL:HG11	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:141:SER:O	1:F:142:LEU:HD23	2.14	0.47
1:C:96:GLY:HA2	1:C:126:GLU:O	2.14	0.47
1:D:137:LEU:O	1:D:275:ARG:NH1	2.47	0.47
1:C:234:LEU:O	1:C:239:MET:HB2	2.14	0.47
1:F:265:THR:HB	1:F:266:PRO:HD2	1.96	0.47
1:G:250:GLN:OE1	1:H:253:VAL:HA	2.14	0.47
1:A:94:ARG:HB3	1:A:142:LEU:HA	1.96	0.47
1:B:283:LEU:O	1:B:287:VAL:HG23	2.15	0.47
1:G:201:GLY:O	1:G:205:TYR:HB3	2.15	0.47
1:C:251:VAL:HG12	1:C:253:VAL:HG23	1.97	0.47
1:D:202:LEU:CD2	1:A:140:GLY:HA3	2.31	0.47
1:E:96:GLY:HA2	1:E:126:GLU:O	2.14	0.47
1:G:147:LEU:HD13	1:G:151:PHE:HE2	1.79	0.46
1:H:183:ILE:HD12	1:H:259:ARG:HB2	1.96	0.46
1:H:200:ILE:CG2	1:H:204:LEU:HB3	2.45	0.46
1:H:282:VAL:O	1:H:282:VAL:HG12	2.14	0.46
1:A:182:GLU:OE2	1:A:262:ALA:HB3	2.15	0.46
1:H:111:ILE:HG23	1:H:123:LEU:CD1	2.45	0.46
1:C:211:SER:HB3	1:C:266:PRO:HG2	1.97	0.46
1:E:97:PHE:O	1:E:127:GLU:HA	2.15	0.46
1:C:265:THR:N	1:C:266:PRO:CD	2.78	0.46
1:A:165:GLU:O	1:A:267:THR:HB	2.16	0.46
1:F:188:PHE:O	1:F:191[B]:GLU:HB2	2.15	0.46
1:H:199:GLU:HG3	1:H:229:ALA:CB	2.44	0.46
1:C:185:LEU:HB2	1:C:263:ASP:OD2	2.14	0.46
1:D:158:HIS:CE1	1:D:276:ARG:NH2	2.84	0.46
1:D:258:TYR:C	1:D:259:ARG:HD2	2.35	0.46
1:A:272:LEU:HD23	1:A:291:PHE:CE1	2.40	0.46
1:G:198:ARG:HD3	1:G:205:TYR:CD1	2.51	0.46
1:C:177:ALA:O	1:C:183:ILE:HD11	2.16	0.46
1:F:115:ARG:HG2	1:F:123:LEU:HD12	1.98	0.46
1:G:267:THR:OG1	1:G:269:LYS:NZ	2.49	0.46
1:F:200:ILE:HG23	1:F:204:LEU:HD23	1.97	0.46
1:F:202:LEU:O	1:F:202:LEU:HD23	2.16	0.45
1:E:144:VAL:HG23	1:E:274:TYR:HB3	1.97	0.45
1:F:231:VAL:O	1:F:235:VAL:HG23	2.17	0.45
1:H:283:LEU:HD12	1:H:283:LEU:O	2.16	0.45
1:E:112:ARG:CD	1:F:238:GLU:HG3	2.45	0.45
1:C:185:LEU:HG	1:C:261:ILE:CG2	2.46	0.45
1:B:196:PHE:HB3	1:B:198:ARG:NH2	2.32	0.45
1:H:195:LEU:HB3	1:H:205:TYR:HE1	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:ASP:OD2	1:H:282:VAL:CG2	2.62	0.45
1:A:118:TYR:CZ	1:A:285:ASN:HB3	2.52	0.45
1:C:251:VAL:O	1:C:258:TYR:OH	2.27	0.45
1:A:95:VAL:HG22	1:A:144:VAL:HG12	1.99	0.45
1:F:132:ARG:HG2	1:F:132:ARG:NH2	2.31	0.45
1:D:132:ARG:HG2	1:D:132:ARG:NH2	2.32	0.45
1:E:171:MET:CE	1:E:259:ARG:HG2	2.47	0.44
1:G:96:GLY:HA2	1:G:126:GLU:O	2.17	0.44
1:H:195:LEU:HB3	1:H:205:TYR:CE1	2.52	0.44
1:F:200:ILE:HD11	1:F:244:VAL:HG12	1.99	0.44
1:G:112:ARG:HG3	1:H:238:GLU:OE1	2.17	0.44
1:E:89:GLU:O	1:E:90:LEU:HD23	2.18	0.44
1:F:205:TYR:OH	1:G:94:ARG:NH2	2.48	0.44
1:C:114:PHE:CD1	1:C:286:PHE:HA	2.52	0.44
1:E:125:LEU:HD11	1:F:239:MET:SD	2.58	0.44
1:F:189:ARG:HA	1:F:217:PHE:CD2	2.52	0.44
1:E:219:PRO:HG2	1:E:221:ILE:HD11	1.99	0.44
1:G:205:TYR:CE1	1:G:209:ILE:HD12	2.53	0.44
1:A:162:LEU:HD12	1:A:162:LEU:HA	1.79	0.44
1:B:107:VAL:HB	1:B:108:PRO:HD3	1.98	0.44
1:F:195:LEU:HD21	1:F:208:VAL:HG12	2.00	0.44
1:F:244:VAL:HB	1:F:248:MET:HG3	2.00	0.44
1:E:162:LEU:HA	1:E:162:LEU:HD12	1.79	0.44
1:F:263:ASP:O	1:F:265:THR:HG23	2.17	0.44
1:D:158:HIS:O	1:D:273:ALA:HA	2.18	0.44
1:G:275:ARG:NH2	1:G:278:ASP:OD2	2.50	0.44
1:D:148:ARG:NH1	1:D:269:LYS:O	2.51	0.44
1:A:137:LEU:HD23	1:A:137:LEU:HA	1.83	0.44
1:E:204:LEU:O	1:E:208:VAL:HG23	2.18	0.44
1:G:228:ILE:HA	1:G:231:VAL:HG23	2.00	0.43
1:A:110:ALA:HB1	1:A:290:VAL:HG22	2.00	0.43
1:F:169:ILE:HD13	1:F:188:PHE:CZ	2.52	0.43
1:G:159:LEU:HD12	1:G:273:ALA:HB2	1.99	0.43
1:D:200:ILE:HG22	1:D:204:LEU:HB3	1.99	0.43
1:F:182:GLU:HG2	1:F:262:ALA:HB2	1.99	0.43
1:F:237:ALA:O	1:F:238:GLU:HB2	2.18	0.43
1:H:232:ILE:O	1:H:235:VAL:HB	2.17	0.43
1:A:103:PHE:HB2	1:A:248:MET:HE2	2.01	0.43
1:G:159:LEU:HD12	1:G:159:LEU:HA	1.84	0.43
1:H:110:ALA:HB1	1:H:290:VAL:HG22	1.99	0.43
1:C:127:GLU:HB3	1:D:228:ILE:CG2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:228:ILE:HG22	1:H:230:SER:H	1.83	0.43
1:E:159:LEU:HA	1:E:159:LEU:HD12	1.60	0.43
1:F:144:VAL:HB	1:F:283:LEU:HD13	2.00	0.43
1:C:194:LEU:HD23	1:C:194:LEU:HA	1.80	0.43
1:H:169:ILE:HD13	1:H:169:ILE:HG21	1.82	0.43
1:B:93:LEU:HB3	1:B:123:LEU:HD23	2.00	0.43
1:E:251:VAL:HG12	1:E:253:VAL:HG23	2.00	0.43
1:C:235:VAL:HG13	1:C:240:GLY:O	2.19	0.43
1:A:161:MET:HG3	1:A:162:LEU:N	2.33	0.43
1:A:251:VAL:HG12	1:A:253:VAL:HG23	2.01	0.43
1:G:171:MET:HE2	1:G:178:ALA:N	2.34	0.43
1:C:92:VAL:HG23	1:C:122:ARG:O	2.18	0.42
1:D:194:LEU:HD23	1:D:194:LEU:HA	1.66	0.42
1:B:198:ARG:CD	1:B:198:ARG:H	2.32	0.42
1:H:168:MET:HE3	1:H:249:SER:HB3	2.01	0.42
1:F:104:ASN:OD1	1:F:106:VAL:HG13	2.18	0.42
1:C:90:LEU:HD23	1:C:91:GLY:N	2.34	0.42
1:B:194:LEU:HD13	1:B:234:LEU:HB2	2.00	0.42
1:E:109:THR:HG22	1:E:112:ARG:HH21	1.84	0.42
1:C:127:GLU:O	1:C:127:GLU:HG2	2.15	0.42
1:F:181:GLU:CD	1:F:181:GLU:H	2.23	0.42
1:A:233:ASN:HB3	1:B:102:ALA:HB1	2.01	0.42
1:F:89:GLU:O	1:F:120:ASP:HB2	2.20	0.42
1:H:288:LEU:HD23	1:H:288:LEU:HA	1.87	0.42
1:E:227:GLN:O	1:E:230:SER:N	2.47	0.42
1:F:261:ILE:HG21	1:F:265:THR:CG2	2.50	0.42
1:D:185:LEU:HD23	1:D:185:LEU:HA	1.69	0.42
1:D:200:ILE:HG23	1:D:204:LEU:HD23	2.02	0.42
1:F:204:LEU:O	1:F:208:VAL:HG23	2.19	0.42
1:A:98:THR:N	1:A:101:SER:OG	2.53	0.42
1:G:148:ARG:HG2	1:G:148:ARG:HH11	1.84	0.42
1:H:267:THR:OG1	1:H:269:LYS:HE3	2.19	0.42
1:B:185:LEU:HA	1:B:185:LEU:HD23	1.88	0.42
1:H:213:ARG:NH2	1:H:218:GLU:OE2	2.47	0.42
1:B:170:VAL:HG22	1:B:235:VAL:HG11	2.01	0.42
1:F:94:ARG:HB3	1:F:142:LEU:HA	2.02	0.42
1:H:259:ARG:HG3	1:H:259:ARG:HH11	1.85	0.42
1:D:167:MET:HG2	1:D:245:PRO:HA	2.02	0.41
1:G:252:LYS:O	1:H:250:GLN:HG2	2.19	0.41
1:A:103:PHE:HB2	1:A:248:MET:CE	2.51	0.41
1:A:151:PHE:H	1:A:151:PHE:HD1	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:95:VAL:HG21	1:F:123:LEU:HD22	2.01	0.41
1:F:95:VAL:O	1:F:125:LEU:HA	2.20	0.41
1:C:221:ILE:HD13	1:C:221:ILE:HA	1.74	0.41
1:B:108:PRO:O	1:B:112:ARG:HB2	2.20	0.41
1:F:202:LEU:HD23	1:F:202:LEU:C	2.40	0.41
1:F:272:LEU:HD23	1:F:291:PHE:HE2	1.85	0.41
1:A:234:LEU:HD23	1:A:234:LEU:HA	1.86	0.41
1:B:125:LEU:HA	1:B:125:LEU:HD23	1.68	0.41
1:B:258:TYR:O	1:B:259:ARG:HD3	2.19	0.41
1:G:225:ALA:HB3	1:G:231:VAL:HG22	2.02	0.41
1:H:158:HIS:CE1	1:H:276:ARG:HG2	2.54	0.41
1:D:95:VAL:HG13	1:D:144:VAL:HG23	2.01	0.41
1:A:183:ILE:HG23	1:A:188:PHE:CE2	2.55	0.41
1:B:204:LEU:HD21	1:B:245:PRO:HG3	2.02	0.41
1:G:236:ALA:HB2	1:G:256:VAL:HG13	2.01	0.41
1:B:101:SER:O	1:B:104:ASN:HB3	2.20	0.41
1:B:209:ILE:HG23	1:B:219:PRO:HG3	2.02	0.41
1:F:248:MET:C	1:F:250:GLN:H	2.24	0.41
1:F:270:LEU:HD12	1:F:270:LEU:HA	1.69	0.41
1:H:212:CYS:HB3	1:H:217:PHE:CE1	2.56	0.41
1:F:183:ILE:HD12	1:F:183:ILE:N	2.36	0.41
1:C:103:PHE:HA	1:D:233:ASN:ND2	2.35	0.41
1:E:185:LEU:HD23	1:E:185:LEU:HA	1.82	0.41
1:C:116:ARG:O	1:C:119:PRO:HD3	2.21	0.41
1:C:284:ARG:HA	1:C:287:VAL:HB	2.03	0.41
1:D:108:PRO:O	1:D:112:ARG:HB3	2.21	0.41
1:H:142:LEU:HA	1:H:142:LEU:HD23	1.79	0.41
1:B:128:ASP:OD1	1:B:129:ASN:N	2.48	0.41
1:B:165:GLU:OE1	1:B:247:SER:OG	2.30	0.40
1:E:258:TYR:O	1:E:259:ARG:HD3	2.21	0.40
1:H:282:VAL:O	1:H:282:VAL:CG1	2.68	0.40
1:C:148:ARG:NE	1:C:204:LEU:HD11	2.36	0.40
1:C:188:PHE:CD2	1:C:191:GLU:HG3	2.56	0.40
1:D:177:ALA:HB1	1:D:183:ILE:HD13	2.01	0.40
1:E:175:HIS:CG	1:E:176:PRO:HD2	2.56	0.40
1:G:86:VAL:N	1:G:89:GLU:OE2	2.54	0.40
1:C:124:GLN:HG3	1:B:224:LEU:CD1	2.51	0.40
1:D:118:TYR:CZ	1:D:285:ASN:ND2	2.89	0.40
1:D:205:TYR:OH	1:A:94:ARG:NH2	2.54	0.40
1:E:94:ARG:HG3	1:E:143:ASP:OD2	2.21	0.40
1:G:166:PRO:O	1:G:246:ALA:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:PHE:CZ	1:C:154:SER:HA	2.56	0.40
1:A:98:THR:HG1	1:A:101:SER:CB	2.35	0.40
1:B:103:PHE:CD2	1:B:248:MET:HG2	2.56	0.40
1:E:104:ASN:OD1	1:E:106:VAL:HG22	2.21	0.40
1:F:236:ALA:HB2	1:F:256:VAL:HG13	2.04	0.40
1:G:188:PHE:HD2	1:G:191:GLU:HG2	1.85	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	198/292 (68%)	191 (96%)	7 (4%)	0	100	100
1	B	200/292 (68%)	190 (95%)	9 (4%)	1 (0%)	29	62
1	C	205/292 (70%)	192 (94%)	13 (6%)	0	100	100
1	D	202/292 (69%)	188 (93%)	13 (6%)	1 (0%)	29	62
1	E	205/292 (70%)	195 (95%)	10 (5%)	0	100	100
1	F	204/292 (70%)	193 (95%)	11 (5%)	0	100	100
1	G	205/292 (70%)	190 (93%)	15 (7%)	0	100	100
1	H	205/292 (70%)	193 (94%)	12 (6%)	0	100	100
All	All	1624/2336 (70%)	1532 (94%)	90 (6%)	2 (0%)	51	82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	226	PRO
1	B	226	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	143/238 (60%)	142 (99%)	1 (1%)	84	90
1	B	158/238 (66%)	156 (99%)	2 (1%)	69	82
1	C	156/238 (66%)	152 (97%)	4 (3%)	46	71
1	D	158/238 (66%)	156 (99%)	2 (1%)	69	82
1	E	152/238 (64%)	151 (99%)	1 (1%)	84	90
1	F	166/238 (70%)	165 (99%)	1 (1%)	86	91
1	G	159/238 (67%)	155 (98%)	4 (2%)	47	71
1	H	161/238 (68%)	158 (98%)	3 (2%)	57	76
All	All	1253/1904 (66%)	1235 (99%)	18 (1%)	67	81

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	151	PHE
1	C	156	ARG
1	C	161	MET
1	C	180	TYR
1	D	199	GLU
1	D	274	TYR
1	A	97	PHE
1	B	263	ASP
1	B	284	ARG
1	E	129	ASN
1	F	284	ARG
1	G	127	GLU
1	G	156	ARG
1	G	174	ASN
1	G	249	SER
1	H	147	LEU
1	H	154	SER
1	H	269	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	158	HIS
1	D	223	GLN
1	D	233	ASN
1	D	285	ASN
1	B	175	HIS
1	E	124	GLN
1	E	260	HIS
1	E	285	ASN
1	F	233	ASN
1	H	129	ASN
1	H	227	GLN
1	H	233	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 [i](#)

There are no ligands in this entry.

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	204/292 (69%)	-0.18	2 (0%) 82 82	18, 44, 83, 96	0
1	B	202/292 (69%)	-0.31	0 100 100	19, 38, 60, 83	0
1	C	207/292 (70%)	-0.23	4 (1%) 66 64	21, 44, 76, 100	0
1	D	204/292 (69%)	-0.33	0 100 100	19, 39, 61, 76	0
1	E	207/292 (70%)	-0.11	3 (1%) 75 74	27, 45, 81, 93	0
1	F	205/292 (70%)	-0.24	2 (0%) 82 82	21, 40, 69, 83	0
1	G	207/292 (70%)	-0.21	2 (0%) 82 82	25, 45, 82, 93	0
1	H	207/292 (70%)	-0.19	3 (1%) 75 74	22, 42, 67, 98	0
All	All	1643/2336 (70%)	-0.22	16 (0%) 82 82	18, 42, 77, 100	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	292	PRO	4.7
1	H	182	GLU	4.4
1	H	264	GLN	3.8
1	C	86	VAL	3.6
1	C	179	SER	3.4
1	A	201	GLY	3.4
1	F	180	TYR	3.3
1	H	179	SER	2.9
1	C	181	GLU	2.9
1	G	263	ASP	2.8
1	E	86	VAL	2.7
1	C	265	THR	2.5
1	A	198	ARG	2.4
1	E	197	PRO	2.4
1	F	264	GLN	2.3
1	G	264	GLN	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.