



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:31 PM UTC

PDB ID : 7W08 / pdb_00007w08
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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

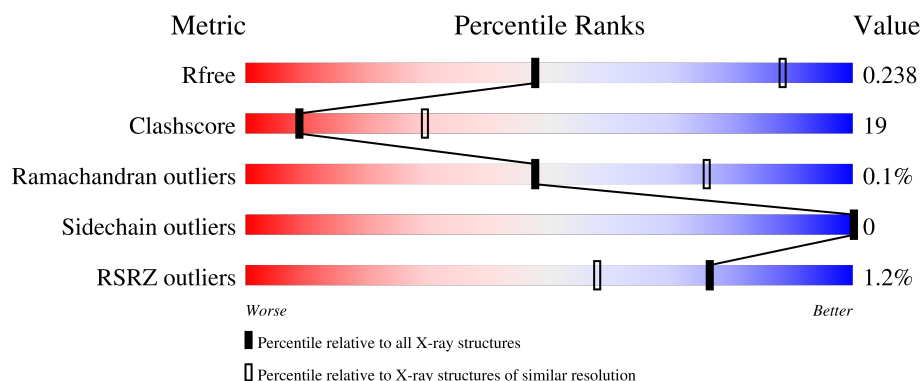
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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}	180053	1605 (3.30-3.22)
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)
RSRZ outliers	180081	1605 (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	% 46% 23% . 30%
1	B	292	45% 23% . 31%
1	C	292	45% 26% 29%
1	D	292	% 48% 21% . 30%
1	E	292	% 48% 23% 29%

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Mol	Chain	Length	Quality of chain
1	F	292	% 43% 26% 30%
1	G	292	% 49% 21% 29%
1	H	292	% 48% 22% 29%

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.

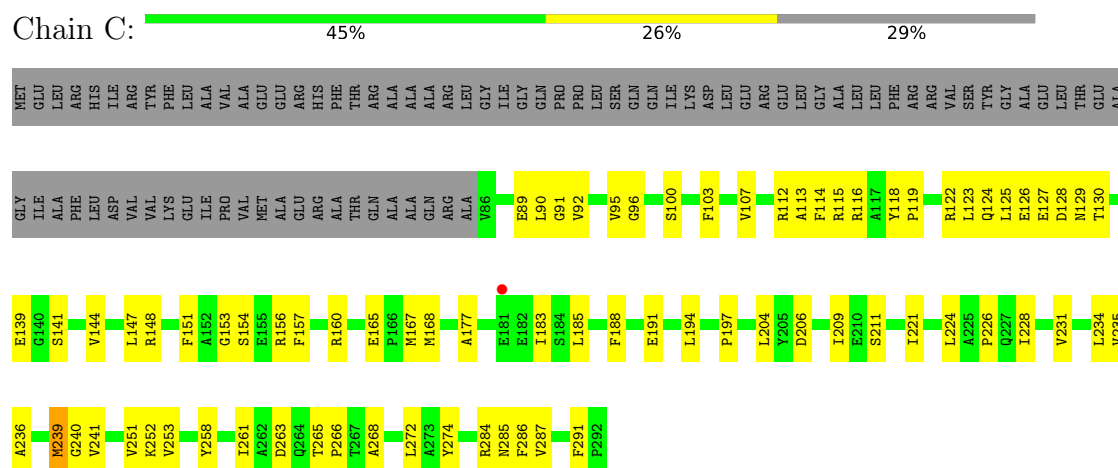
- Molecule 1 is a protein called Transcriptional regulator, LysR family.

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			

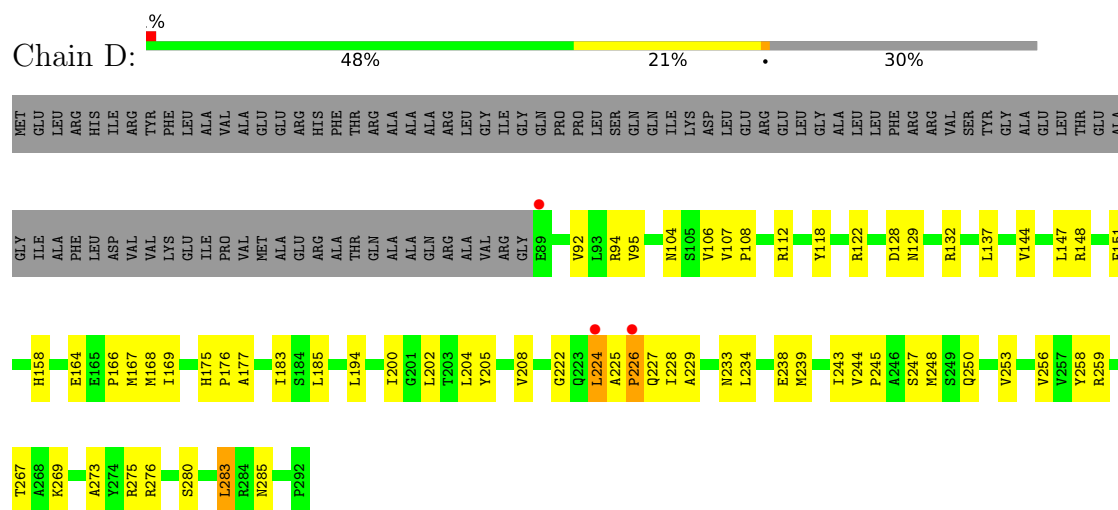
3 Residue-property plots

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.

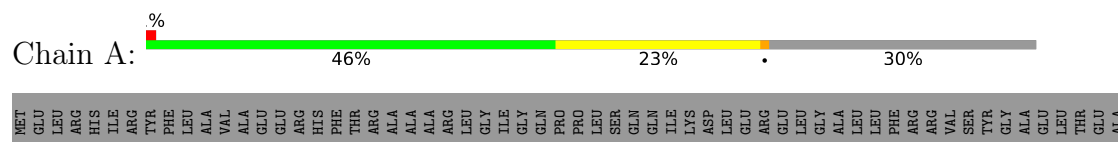
- Molecule 1: Transcriptional regulator, LysR family

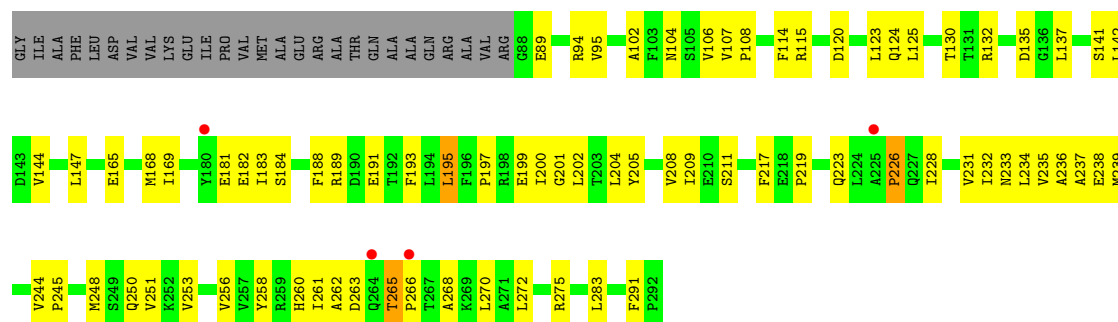


- Molecule 1: Transcriptional regulator, LysR family

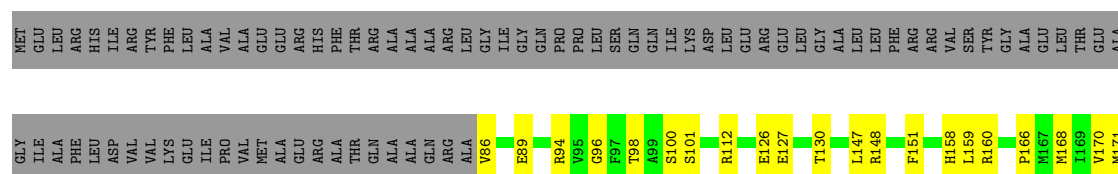


- Molecule 1: Transcriptional regulator, LysR family

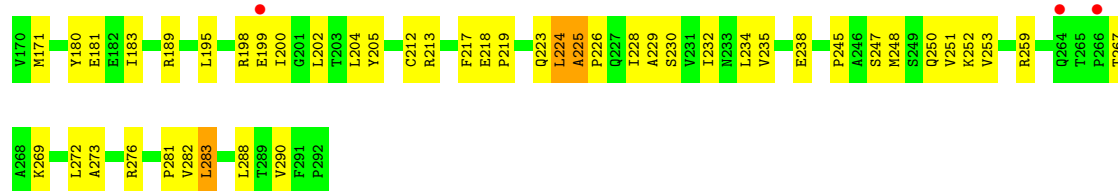
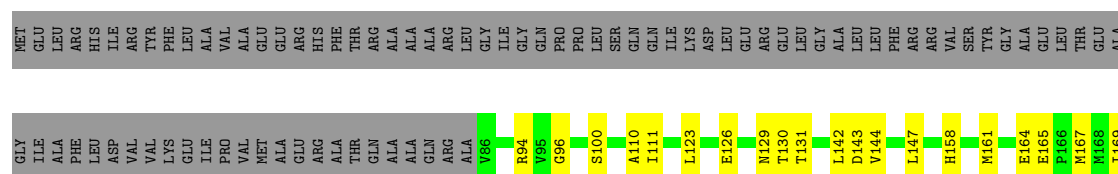




• Molecule 1: Transcriptional regulator, LysR family



• 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 32.96 – 3.25	Depositor EDS
% Data completeness (in resolution range)	91.4 (32.96-3.25) 87.2 (32.96-3.25)	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 0.200 , 0.238	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)	0.34 , 47.1	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	0.64	1/1495 (0.1%)	0.97	3/2038 (0.1%)
1	B	0.73	0/1542	1.09	8/2101 (0.4%)
1	C	0.82	3/1554 (0.2%)	1.12	3/2119 (0.1%)
1	D	0.74	1/1551 (0.1%)	1.08	9/2114 (0.4%)
1	E	0.57	0/1549	0.90	0/2114
1	F	0.80	3/1592 (0.2%)	1.12	5/2167 (0.2%)
1	G	0.82	3/1575 (0.2%)	1.10	7/2145 (0.3%)
1	H	0.81	2/1578 (0.1%)	1.10	6/2149 (0.3%)
All	All	0.75	13/12436 (0.1%)	1.06	41/16947 (0.2%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	245	PRO	N-CA	16.86	1.67	1.47
1	G	244	VAL	C-N	7.90	1.44	1.33
1	H	281	PRO	C-O	-7.02	1.14	1.24
1	F	226	PRO	C-O	-7.00	1.14	1.24
1	F	197	PRO	C-O	-6.94	1.15	1.24
1	F	195	LEU	C-O	-6.40	1.16	1.24
1	H	283	LEU	C-O	-6.32	1.16	1.24
1	G	245	PRO	C-O	-5.72	1.16	1.23
1	D	226	PRO	C-O	-5.46	1.16	1.24
1	C	107	VAL	CA-CB	-5.40	1.51	1.54
1	A	228	ILE	C-O	-5.38	1.17	1.24
1	C	129	ASN	C-O	-5.36	1.17	1.23
1	C	128	ASP	C-O	-5.21	1.18	1.23

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	244	VAL	CA-C-N	15.85	136.06	119.89
1	G	244	VAL	C-N-CA	15.85	136.06	119.89
1	B	226	PRO	N-CA-CB	14.43	118.41	103.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	245	PRO	N-CA-C	-9.10	96.63	111.11
1	H	224	LEU	CB-CA-C	-8.63	99.75	112.09
1	H	223	GLN	CB-CA-C	-8.23	95.58	109.65
1	D	168	MET	CB-CG-SD	-7.70	89.60	112.70
1	H	224	LEU	N-CA-C	-7.58	91.83	107.67
1	H	225	ALA	N-CA-C	-7.38	93.50	109.81
1	C	239	MET	CB-CG-SD	-7.16	91.23	112.70
1	B	227	GLN	CB-CA-C	-6.96	99.40	109.84
1	G	244	VAL	O-C-N	-6.75	115.79	121.40
1	D	226	PRO	N-CA-CB	6.48	110.06	103.25
1	G	245	PRO	CA-N-CD	-6.47	102.94	112.00
1	B	197	PRO	CB-CA-C	-6.41	101.76	111.44
1	D	224	LEU	CA-C-N	6.27	129.07	120.67
1	D	224	LEU	C-N-CA	6.27	129.07	120.67
1	D	222	GLY	CA-C-N	-6.21	114.04	122.86
1	D	222	GLY	C-N-CA	-6.21	114.04	122.86
1	H	225	ALA	CB-CA-C	-6.05	98.25	110.17
1	D	283	LEU	CB-CG-CD1	-5.99	92.72	110.70
1	B	226	PRO	CA-C-N	5.99	129.62	121.05
1	B	226	PRO	C-N-CA	5.99	129.62	121.05
1	F	201	GLY	CA-C-O	-5.66	118.54	122.22
1	A	229	ALA	CA-C-N	5.64	128.41	120.28
1	A	229	ALA	C-N-CA	5.64	128.41	120.28
1	C	206	ASP	CB-CA-C	5.56	120.02	110.79
1	F	223	GLN	CA-C-N	-5.56	114.97	121.64
1	F	223	GLN	C-N-CA	-5.56	114.97	121.64
1	G	246	ALA	CA-C-N	5.55	128.27	120.28
1	G	246	ALA	C-N-CA	5.55	128.27	120.28
1	F	265	THR	OG1-CB-CG2	5.51	120.33	109.30
1	A	116	ARG	CB-CG-CD	5.42	123.77	111.30
1	F	265	THR	N-CA-C	5.42	116.82	110.31
1	B	198	ARG	CB-CG-CD	-5.28	99.17	111.30
1	C	226	PRO	N-CA-C	5.17	120.87	113.47
1	H	225	ALA	N-CA-CB	5.12	119.47	110.37
1	B	196	PHE	CA-CB-CG	5.11	118.91	113.80
1	D	229	ALA	CA-C-N	5.04	131.16	121.54
1	D	229	ALA	C-N-CA	5.04	131.16	121.54
1	B	199	GLU	CB-CA-C	5.04	117.19	109.03

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	54	0
1	B	1512	0	1495	68	0
1	C	1525	0	1493	71	0
1	D	1521	0	1491	59	0
1	E	1518	0	1468	56	0
1	F	1561	0	1550	73	0
1	G	1544	0	1503	53	0
1	H	1548	0	1546	72	0
All	All	12198	0	11937	448	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 19.

All (448) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:G:245:PRO:N	1:G:245:PRO:C	2.40	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:A:168:MET:HE2	1:A:260:HIS:CE1	2.18	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: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	0.74
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:F:199:GLU:OE2	1:F:231:VAL:CG1	2.37	0.68
1:G:228:ILE:HG22	1:G:248:MET:HE2	1.75	0.68
1:H:225:ALA:O	1:H:226:PRO:C	2.36	0.68
1:E:171:MET:HE1	1:E:183:ILE:HD11	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:167:MET:HE3	1:A:243:ILE:HG22	1.78	0.65
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:E:139:GLU:CD	1:H:198:ARG:HH21	2.05	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: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:D:167:MET:HE3	1:D:245:PRO:HD3	1.81	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:G:168:MET:HE1	1:G:249:SER:HB2	1.80	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:A:161:MET:HE1	1:A:269:LYS:HB3	1.80	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
1:F:168:MET:HE2	1:F:260:HIS:CD2	2.36	0.61
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:G:171:MET:HE1	1:G:177:ALA:HB3	1.82	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:H:171:MET:HE1	1:H:183:ILE:HD13	1.85	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: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:E:168:MET:HE3	1:E:260:HIS:CE1	2.40	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: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
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:168:MET:HE3	1:B:246:ALA:HB1	1.87	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:G:168:MET:HE2	1:G:246:ALA:HA	1.87	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:228:ILE:HG22	1:E:248:MET:HE2	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:H:213:ARG:HE	1:H:218:GLU:CD	2.14	0.54
1:E:129:ASN:HB2	1:E:132:ARG:H	1.72	0.54
1:E:171:MET:HE1	1:E:183:ILE:CD1	2.37	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:C:151:PHE:CD1	1:C:151:PHE:C	2.86	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
1:G:127:GLU:CD	1:H:228:ILE:HG23	2.34	0.52
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:G:228:ILE:CG2	1:G:248:MET:HE2	2.39	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:H:167:MET:HE1	1:H:204:LEU:HG	1.93	0.51
1:D:258:TYR:C	1:D:259:ARG:HD2	2.35	0.51
1:B:251:VAL:O	1:B:253:VAL:N	2.41	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:B:167:MET:HE2	1:B:243:ILE:HG22	1.92	0.50
1:A:259:ARG:HG3	1:A:259:ARG:HH11	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:MET:HE2	1:A:260:HIS:NE2	2.27	0.50
1:C:113:ALA:O	1:C:116:ARG:HB3	2.12	0.50
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:D:147:LEU:HD13	1:D:151:PHE:CE2	2.46	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:E:239:MET:HE2	1:F:125:LEU:CD1	2.44	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:161:MET:HE3	1:H:161:MET:HB3	1.70	0.47
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:F:181:GLU:CD	1:F:181:GLU:H	2.23	0.47
1:G:175:HIS:CG	1:G:176:PRO:HD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:F:202:LEU:C	1:F:202:LEU:HD23	2.40	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:103:PHE:HB2	1:A:248:MET:HE1	1.98	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:A:272:LEU:HD23	1:A:291:PHE:CE1	2.40	0.46
1:B:98:THR:O	1:B:101:SER:OG	2.34	0.46
1:F:248:MET:C	1:F:250:GLN:H	2.24	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:HD23	1:F:202:LEU:O	2.16	0.45
1:E:144:VAL:HG23	1:E:274:TYR:HB3	1.97	0.45
1:D:239:MET:HB3	1:D:239:MET:HE2	1.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:G:239:MET:HE2	1:G:239:MET:HB3	1.54	0.45
1:H:195:LEU:HB3	1:H:205:TYR:HE1	1.80	0.45
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:H:248:MET:HE2	1:H:248:MET:HB3	1.72	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:E:204:LEU:O	1:E:207:SER:OG	2.36	0.43
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:G:171:MET:HE3	1:G:175:HIS:ND1	2.32	0.43
1:H:171:MET:HE1	1:H:183:ILE:CD1	2.47	0.43
1:H:232:ILE:O	1:H:235:VAL:HB	2.17	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
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:G:171:MET:HE3	1:G:175:HIS:HB3	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: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: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:A:100:SER:HB2	1:A:248:MET:HE3	2.01	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:124:GLN:CD	1:B:225:ALA:HB2	2.44	0.42
1:C:127:GLU:O	1:C:127:GLU:HG2	2.15	0.42
1:D:248:MET:HE3	1:D:248:MET:HB3	1.72	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:F:248:MET:HB3	1:F:248:MET:HE3	1.70	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
1:F:95:VAL:O	1:F:125:LEU:HA	2.20	0.41
1:F:95:VAL:HG21	1:F:123:LEU:HD22	2.01	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: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:B:207:SER:O	1:B:211:SER:OG	2.38	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:114:PHE:CD1	1:F:114:PHE:C	2.98	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:B:161:MET:HE1	1:B:269:LYS:HG2	2.02	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:168:MET:HE2	1:C:258:TYR:CB	2.51	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:E:239:MET:HE2	1:F:125:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
1:C:151:PHE:CZ	1:C:154:SER:HA	2.56	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%)	24	55
1	C	205/292 (70%)	192 (94%)	13 (6%)	0	100	100
1	D	202/292 (69%)	188 (93%)	13 (6%)	1 (0%)	24	55
1	E	205/292 (70%)	195 (95%)	10 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
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%)	48	76

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%)	143 (100%)	0	100	100
1	B	158/238 (66%)	158 (100%)	0	100	100
1	C	156/238 (66%)	156 (100%)	0	100	100
1	D	158/238 (66%)	158 (100%)	0	100	100
1	E	152/238 (64%)	152 (100%)	0	100	100
1	F	166/238 (70%)	166 (100%)	0	100	100
1	G	159/238 (67%)	159 (100%)	0	100	100
1	H	161/238 (68%)	161 (100%)	0	100	100
All	All	1253/1904 (66%)	1253 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	129	ASN

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Mol	Chain	Res	Type
1	D	124	GLN
1	D	158	HIS
1	D	223	GLN
1	D	233	ASN
1	D	285	ASN
1	A	104	ASN
1	A	260	HIS
1	B	175	HIS
1	B	233	ASN
1	E	124	GLN
1	E	174	ASN
1	E	233	ASN
1	E	285	ASN
1	F	233	ASN
1	H	129	ASN
1	H	227	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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 ⓘ

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%)	0.09	4 (1%) 65 46	18, 44, 83, 96	0
1	B	202/292 (69%)	-0.17	1 (0%) 87 76	19, 38, 60, 83	0
1	C	207/292 (70%)	-0.05	1 (0%) 87 76	21, 44, 76, 100	0
1	D	204/292 (69%)	-0.16	3 (1%) 72 53	19, 39, 61, 76	0
1	E	207/292 (70%)	0.00	2 (0%) 79 62	27, 45, 81, 93	0
1	F	205/292 (70%)	-0.11	4 (1%) 65 46	21, 40, 69, 83	1 (0%)
1	G	207/292 (70%)	0.08	2 (0%) 79 62	25, 45, 82, 93	0
1	H	207/292 (70%)	-0.02	3 (1%) 73 54	22, 42, 67, 98	0
All	All	1643/2336 (70%)	-0.04	20 (1%) 76 58	18, 42, 77, 100	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	225	ALA	4.0
1	H	264	GLN	3.2
1	A	226	PRO	3.0
1	E	181	GLU	2.9
1	E	292	PRO	2.9
1	A	202	LEU	2.9
1	D	226	PRO	2.8
1	A	129	ASN	2.6
1	G	200	ILE	2.5
1	F	266	PRO	2.4
1	F	180	TYR	2.4
1	G	263	ASP	2.4
1	B	226	PRO	2.2
1	H	266	PRO	2.2
1	D	224	LEU	2.2
1	D	89	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	222	GLY	2.2
1	F	264	GLN	2.1
1	C	181	GLU	2.1
1	H	199	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.