



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 04:04 PM BST

PDB ID : 6UON
Title : Molecular basis for tumor infiltrating TCR recognition of hotspot KRAS-G12D mutation
Authors : Sun, P.D.; Sim, M.J.W.
Deposited on : 2019-10-15
Resolution : 3.50 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 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	:	2.13.1
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)	:	2.13.1

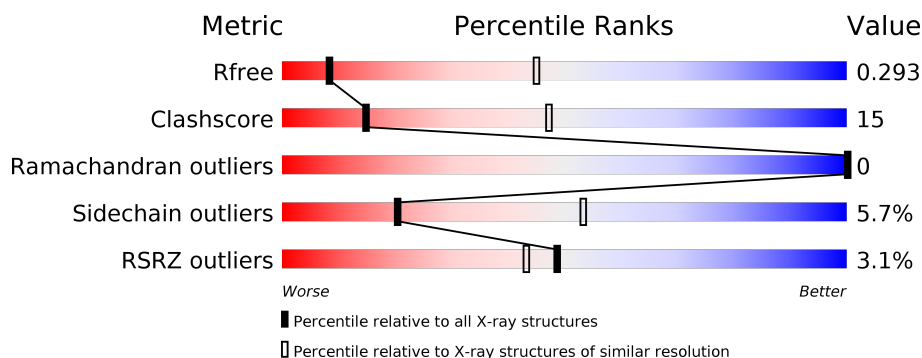
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	G	202	<div> <div>4%</div> <div> <div></div> <div>59%</div> <div>31%</div> <div>9%</div> <div>.</div> </div> </div>
1	I	202	<div> <div>8%</div> <div> <div></div> <div>59%</div> <div>33%</div> <div>.</div> <div>.</div> </div> </div>
2	H	241	<div> <div>5%</div> <div> <div></div> <div>72%</div> <div>24%</div> <div>..</div> </div> </div>
2	J	241	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>22%</div> <div>..</div> </div> </div>
3	A	274	<div> <div>3%</div> <div> <div></div> <div>77%</div> <div>21%</div> <div>.</div> </div> </div>
3	D	274	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>20%</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	B	99	<div><div><div>%</div><div><div></div><div>87%</div><div>11%</div><div>..</div></div></div></div>
4	E	99	<div><div><div>%</div><div><div></div><div>78%</div><div>20%</div><div>.</div></div></div></div>
5	C	10	<div><div><div></div><div><div>80%</div><div>20%</div></div></div></div>
5	F	10	<div><div><div></div><div><div>60%</div><div>30%</div><div>10%</div></div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13102 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 TCR-V-alpha-12-02*01.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	199	Total	C	N	O	S	0	0	0
			1551	967	256	319	9			
1	I	193	Total	C	N	O	S	0	0	0
			1507	943	249	306	9			

- Molecule 2 is a protein called TCR-V-beta-10-2*01.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	238	Total	C	N	O	S	0	0	0
			1904	1201	332	364	7			
2	J	238	Total	C	N	O	S	0	0	0
			1904	1201	332	364	7			

- Molecule 3 is a protein called HLA class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	273	Total	C	N	O	S	0	0	0
			2241	1395	414	426	6			
3	D	273	Total	C	N	O	S	0	0	0
			2241	1395	414	426	6			

- Molecule 4 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	98	Total	C	N	O	S	0	0	0
			820	523	139	156	2			
4	E	97	Total	C	N	O	S	0	0	0
			812	517	138	155	2			

- Molecule 5 is a protein called GLY-ALA-ASP-GLY-VAL-GLY-LYS-SER-ALA-LEU.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	C	10	Total	C	N	O	0	0	0
			61	36	11	14			
5	F	10	Total	C	N	O	0	0	0
			61	36	11	14			

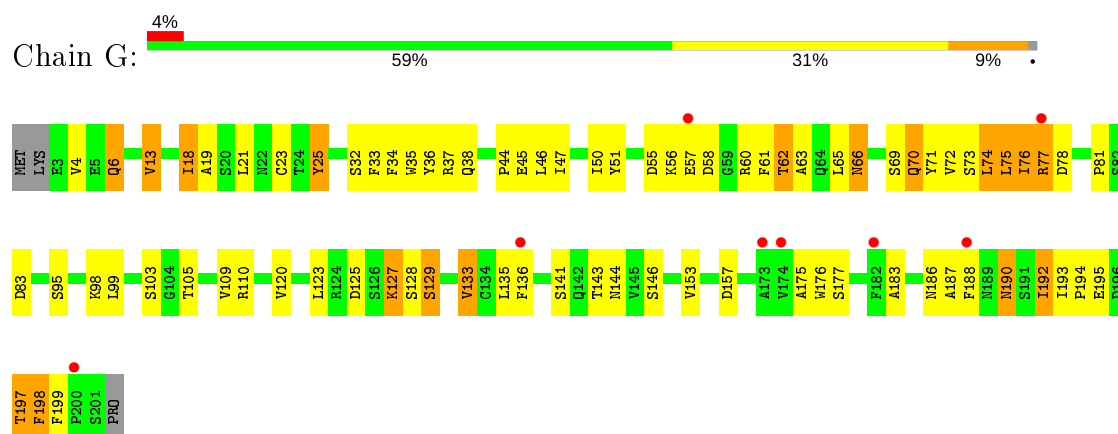
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	3	ASP	GLY	conflict	UNP P01111
F	3	ASP	GLY	conflict	UNP P01111

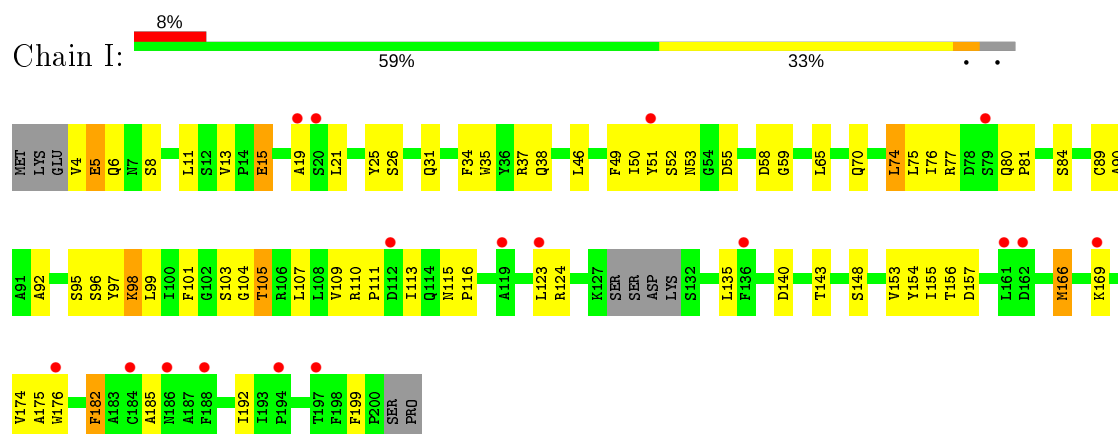
3 Residue-property plots [i](#)

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.

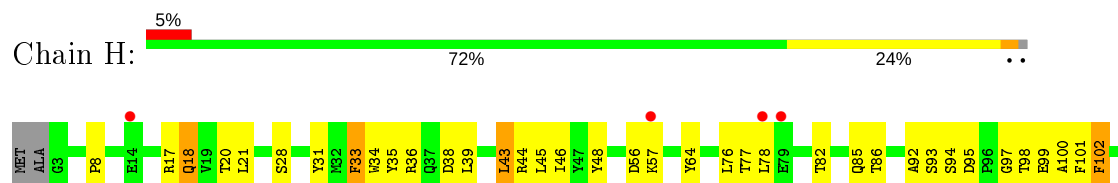
- Molecule 1: TCR-V-alpha-12-02*01

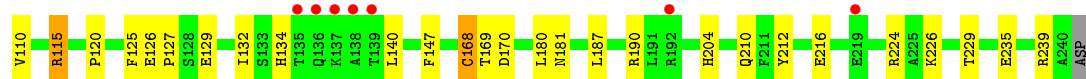


- Molecule 1: TCR-V-alpha-12-02*01

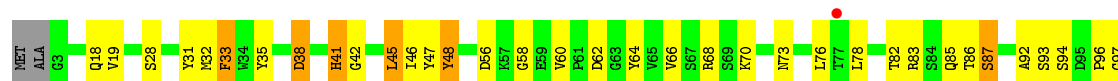


- Molecule 2: TCR-V-beta-10-2*01

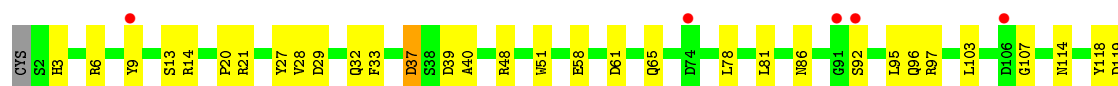
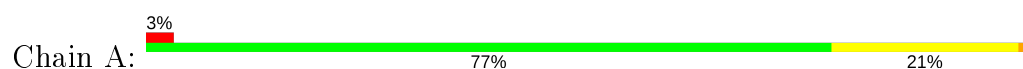




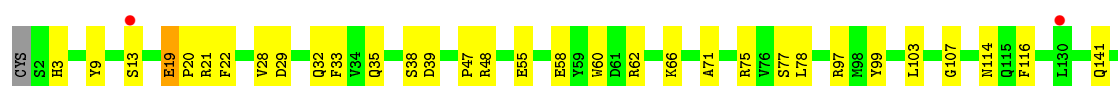
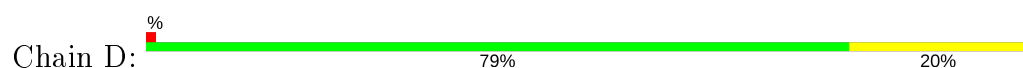
• Molecule 2: TCR-V-beta-10-2*01



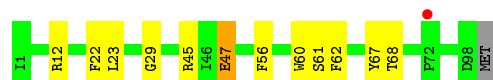
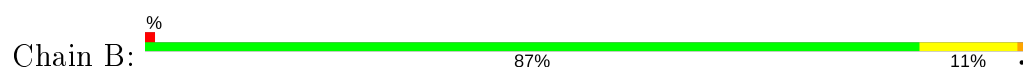
• Molecule 3: HLA class I antigen



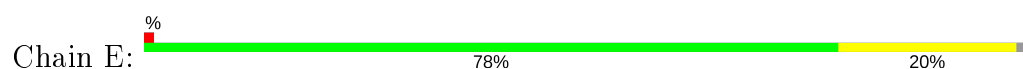
• Molecule 3: HLA class I antigen




• Molecule 4: Beta-2-microglobulin



• Molecule 4: Beta-2-microglobulin



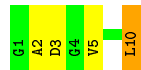
- Molecule 5: GLY-ALA-ASP-GLY-VAL-GLY-LYS-SER-ALA-LEU

Chain C:  80% 20%



- Molecule 5: GLY-ALA-ASP-GLY-VAL-GLY-LYS-SER-ALA-LEU

Chain F:  60% 30% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.40 Å 79.29 Å 197.71 Å 90.00° 91.11° 90.00°	Depositor
Resolution (Å)	49.42 – 3.50 49.42 – 3.46	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.42-3.50) 98.9 (49.42-3.46)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 3.48 Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.252 , 0.294 0.251 , 0.293	Depositor DCC
R_{free} test set	1299 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	112.7	Xtriage
Anisotropy	0.527	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 67.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.048 for h,-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13102	wwPDB-VP
Average B, all atoms (Å ²)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	G	0.42	0/1584	0.68	0/2145
1	I	0.39	0/1539	0.62	0/2084
2	H	0.30	0/1958	0.53	0/2669
2	J	0.28	0/1958	0.50	0/2669
3	A	0.26	0/2303	0.46	0/3125
3	D	0.27	0/2303	0.50	0/3125
4	B	0.24	0/843	0.44	0/1142
4	E	0.24	0/835	0.45	0/1131
5	C	0.51	0/60	0.66	0/77
5	F	0.53	0/60	0.99	0/77
All	All	0.31	0/13443	0.53	0/18244

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	G	1551	0	1461	119	0
1	I	1507	0	1422	66	0
2	H	1904	0	1801	53	0
2	J	1904	0	1801	47	0
3	A	2241	0	2098	52	0
3	D	2241	0	2098	50	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	820	0	785	7	0
4	E	812	0	771	16	0
5	C	61	0	63	5	0
5	F	61	0	63	9	0
All	All	13102	0	12363	386	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 15.

All (386) 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:I:153:VAL:HG21	1:I:182:PHE:CD1	1.66	1.31
1:G:23:CYS:SG	1:G:72:VAL:CG1	2.24	1.25
3:D:77:SER:HB3	5:F:10:LEU:CD1	1.67	1.25
1:G:23:CYS:SG	1:G:72:VAL:HG11	1.78	1.22
1:G:25:TYR:CZ	1:G:70:GLN:HB2	1.80	1.17
1:G:19:ALA:HB3	1:G:76:ILE:HD11	1.16	1.16
1:G:19:ALA:CB	1:G:76:ILE:HD11	1.83	1.09
1:G:81:PRO:HA	1:G:109:VAL:HG11	1.17	1.07
3:D:77:SER:HB3	5:F:10:LEU:HD13	1.34	1.07
1:I:153:VAL:CG2	1:I:182:PHE:CD1	2.42	1.02
1:G:133:VAL:HG12	1:G:176:TRP:CB	1.88	1.02
1:G:19:ALA:HB3	1:G:76:ILE:CD1	1.88	1.02
3:D:13:SER:HB3	3:D:78:LEU:HD13	1.39	1.02
3:D:77:SER:HB3	5:F:10:LEU:HD12	1.39	1.01
2:H:35:TYR:HB3	2:H:43:LEU:HD11	1.41	1.01
1:I:4:VAL:O	1:I:26:SER:HB2	1.62	1.00
1:I:5:GLU:OE2	1:I:25:TYR:HA	1.64	0.98
1:G:133:VAL:CG1	1:G:176:TRP:CB	2.42	0.96
1:G:76:ILE:HD13	1:G:76:ILE:H	1.28	0.96
1:G:183:ALA:HB3	1:G:186:ASN:HB3	1.47	0.96
1:G:25:TYR:OH	1:G:70:GLN:HB2	1.65	0.96
3:D:77:SER:CB	5:F:10:LEU:HD12	1.96	0.94
3:A:32:GLN:OE1	3:A:48:ARG:HG3	1.68	0.94
1:I:153:VAL:HG21	1:I:182:PHE:CE1	2.02	0.94
1:I:21:LEU:HD22	1:I:105:THR:HG21	1.49	0.94
1:G:133:VAL:HG12	1:G:176:TRP:HB3	1.50	0.92
1:G:23:CYS:SG	1:G:72:VAL:HG12	2.09	0.92
1:G:133:VAL:CG1	1:G:176:TRP:HB2	2.00	0.92
1:I:35:TRP:CG	1:I:74:LEU:HD12	2.06	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:133:VAL:CG1	1:G:176:TRP:HB3	2.03	0.89
1:I:15:GLU:HG3	1:I:111:PRO:HA	1.56	0.87
1:I:35:TRP:CG	1:I:74:LEU:CD1	2.59	0.85
1:G:75:LEU:HD23	1:G:75:LEU:O	1.76	0.84
1:G:21:LEU:HD12	1:G:105:THR:HG21	1.62	0.81
1:I:153:VAL:HG21	1:I:182:PHE:HD1	1.40	0.81
1:G:19:ALA:CA	1:G:76:ILE:HD11	2.11	0.80
1:I:21:LEU:HD22	1:I:105:THR:CG2	2.12	0.80
1:G:19:ALA:O	1:G:76:ILE:CD1	2.29	0.80
1:G:105:THR:HG23	1:G:105:THR:O	1.80	0.79
1:G:193:ILE:O	1:G:193:ILE:HD12	1.82	0.79
1:G:19:ALA:O	1:G:76:ILE:HD13	1.81	0.79
1:G:33:PHE:CD1	1:G:65:LEU:HD21	2.18	0.78
1:I:192:ILE:HG22	1:I:192:ILE:O	1.83	0.78
3:A:81:LEU:HD11	5:C:10:LEU:HD11	1.65	0.77
4:E:54:LEU:HD11	4:E:62:PHE:HD2	1.47	0.77
1:G:38:GLN:HB2	1:G:44:PRO:HG3	1.67	0.76
1:I:35:TRP:CD1	1:I:74:LEU:HD12	2.20	0.76
1:G:133:VAL:HG13	1:G:176:TRP:HB2	1.64	0.76
3:D:77:SER:CB	5:F:10:LEU:CD1	2.53	0.75
1:G:183:ALA:O	1:G:187:ALA:N	2.20	0.75
1:G:192:ILE:HD13	1:G:192:ILE:N	2.02	0.75
2:H:35:TYR:HB3	2:H:43:LEU:CD1	2.17	0.74
1:G:81:PRO:CA	1:G:109:VAL:HG11	2.09	0.73
1:I:5:GLU:OE2	1:I:5:GLU:HA	1.88	0.73
1:G:57:GLU:HA	1:G:62:THR:HG23	1.70	0.73
2:H:94:SER:HB3	2:H:100:ALA:HA	1.71	0.73
1:I:153:VAL:CG2	1:I:182:PHE:HD1	1.97	0.72
1:I:98:LYS:O	1:I:98:LYS:HG3	1.90	0.72
3:D:32:GLN:NE2	4:E:53:ASP:OD2	2.21	0.71
2:J:86:THR:HG23	2:J:109:THR:HA	1.73	0.71
3:D:107:GLY:O	3:D:169:ARG:NH1	2.24	0.70
1:G:192:ILE:H	1:G:192:ILE:HD13	1.55	0.70
2:J:97:GLY:HA3	5:C:5:VAL:HG23	1.73	0.70
3:A:28:VAL:HG21	3:A:51:TRP:HH2	1.56	0.70
1:I:84:SER:HB3	1:I:109:VAL:HG12	1.72	0.70
2:J:64:TYR:HD1	2:J:76:LEU:HD11	1.57	0.70
1:G:195:GLU:OE2	1:G:195:GLU:HA	1.90	0.69
3:D:71:ALA:O	3:D:75:ARG:HG3	1.93	0.69
2:J:46:ILE:HD12	2:J:60:VAL:O	1.93	0.69
3:A:28:VAL:HG21	3:A:51:TRP:CH2	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:32:MET:SD	2:J:68:ARG:NH2	2.64	0.69
2:J:46:ILE:HG23	2:J:47:TYR:H	1.57	0.69
2:J:19:VAL:HG13	2:J:78:LEU:HD21	1.74	0.69
1:I:6:GLN:OE1	1:I:6:GLN:N	2.25	0.68
2:H:43:LEU:HD12	2:H:43:LEU:C	2.14	0.68
1:I:15:GLU:HB2	1:I:110:ARG:O	1.94	0.68
2:H:86:THR:HG22	2:H:110:VAL:H	1.57	0.68
1:I:35:TRP:CB	1:I:74:LEU:HD11	2.22	0.68
2:H:224:ARG:HA	2:J:70:LYS:HE3	1.75	0.68
2:J:46:ILE:CD1	2:J:60:VAL:O	2.42	0.67
2:H:180:LEU:O	2:H:180:LEU:HD23	1.94	0.67
1:G:183:ALA:O	1:G:187:ALA:HB2	1.94	0.67
3:A:208:PHE:CE1	3:A:241:PHE:O	2.48	0.67
1:G:21:LEU:CD1	1:G:105:THR:HG21	2.24	0.67
3:D:232:GLU:OE1	4:E:6:LYS:NZ	2.25	0.67
1:I:35:TRP:CB	1:I:74:LEU:CD1	2.73	0.66
2:J:66:VAL:HG12	2:J:76:LEU:HA	1.77	0.66
3:D:35:GLN:NE2	3:D:48:ARG:HH21	1.93	0.66
2:H:38:ASP:OD2	2:H:38:ASP:N	2.29	0.66
3:A:13:SER:HB3	3:A:78:LEU:HD13	1.78	0.66
3:A:208:PHE:H	3:A:208:PHE:HD1	1.43	0.66
3:D:103:LEU:HD21	3:D:165:VAL:HG13	1.78	0.66
1:G:76:ILE:N	1:G:76:ILE:HD13	2.08	0.65
3:D:19:GLU:HG3	3:D:75:ARG:HH22	1.61	0.65
3:D:219:ARG:HG2	3:D:257:TYR:CE2	2.32	0.65
1:G:190:ASN:N	1:G:190:ASN:OD1	2.30	0.65
1:I:103:SER:O	2:J:42:GLY:HA3	1.97	0.65
1:G:76:ILE:CD1	1:G:76:ILE:H	2.05	0.65
1:I:35:TRP:HB2	1:I:74:LEU:HD11	1.78	0.65
1:G:198:PHE:CE2	2:H:134:HIS:CD2	2.86	0.65
1:G:21:LEU:HD21	1:G:74:LEU:HD21	1.78	0.65
3:D:156:ARG:NE	5:F:3:ASP:OD2	2.31	0.64
2:H:35:TYR:CB	2:H:43:LEU:HD11	2.22	0.64
1:I:46:LEU:HD21	1:I:49:PHE:HD1	1.62	0.64
2:J:32:MET:HG2	2:J:93:SER:HB2	1.80	0.64
3:D:77:SER:OG	5:F:10:LEU:HD12	1.98	0.64
1:G:127:LYS:HE3	1:G:127:LYS:N	2.13	0.64
4:B:29:GLY:HA2	4:B:61:SER:HB2	1.79	0.63
2:H:64:TYR:HD1	2:H:76:LEU:HD11	1.63	0.63
1:G:46:LEU:HD22	2:H:99:GLU:HB2	1.80	0.63
3:A:81:LEU:HD13	3:A:118:TYR:HD1	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:81:PRO:HA	1:I:109:VAL:HG11	1.79	0.63
1:G:74:LEU:O	1:G:74:LEU:HG	1.99	0.63
1:G:6:GLN:NE2	1:G:103:SER:O	2.32	0.62
1:G:75:LEU:C	1:G:75:LEU:HD23	2.18	0.62
3:D:19:GLU:HB2	3:D:75:ARG:NH2	2.14	0.62
1:I:31:GLN:N	1:I:31:GLN:OE1	2.32	0.62
3:D:19:GLU:HG3	3:D:75:ARG:NH2	2.15	0.62
1:G:105:THR:O	1:G:105:THR:CG2	2.47	0.62
3:A:208:PHE:HE1	3:A:241:PHE:O	1.83	0.62
3:D:3:HIS:ND1	3:D:29:ASP:OD1	2.28	0.61
3:A:97:ARG:NH1	3:A:114:ASN:OD1	2.33	0.61
1:I:90:ALA:HB2	1:I:101:PHE:CD1	2.35	0.61
3:D:55:GLU:OE2	3:D:55:GLU:HA	1.99	0.61
2:J:38:ASP:OD1	2:J:87:SER:HB3	2.00	0.61
3:A:231:VAL:O	3:A:243:LYS:NZ	2.32	0.61
2:H:120:PRO:HB3	2:H:147:PHE:HB3	1.83	0.61
2:J:83:ARG:HH21	2:J:181:ASN:HD22	1.47	0.61
2:J:46:ILE:HG23	2:J:47:TYR:N	2.14	0.61
3:A:3:HIS:ND1	3:A:29:ASP:OD2	2.33	0.61
1:G:123:LEU:CD1	1:G:135:LEU:HD23	2.31	0.60
1:G:143:THR:HG21	1:G:193:ILE:HG22	1.83	0.60
2:H:126:GLU:OE1	2:H:239:ARG:NH2	2.34	0.60
2:H:36:ARG:NH1	2:H:64:TYR:OH	2.34	0.60
3:D:35:GLN:HE21	3:D:48:ARG:HH21	1.49	0.60
1:G:63:ALA:HA	1:G:73:SER:O	2.01	0.60
3:A:208:PHE:HE1	3:A:241:PHE:C	2.04	0.60
1:G:32:SER:HA	1:G:51:TYR:HD1	1.66	0.60
1:G:50:ILE:HG21	1:G:65:LEU:HD13	1.83	0.60
1:G:120:VAL:HG22	1:G:136:PHE:HB2	1.84	0.59
3:A:14:ARG:HH11	3:A:21:ARG:HB2	1.66	0.59
4:E:70:PHE:HE2	4:E:72:PRO:HB3	1.66	0.59
2:H:8:PRO:HG2	2:H:21:LEU:HD22	1.84	0.59
3:A:224:GLN:O	3:A:228:THR:OG1	2.20	0.59
2:J:82:THR:N	2:J:85:GLN:OE1	2.36	0.58
2:H:216:GLU:N	2:H:216:GLU:OE2	2.36	0.58
1:G:133:VAL:HG21	2:H:125:PHE:HD1	1.68	0.58
1:G:193:ILE:HD12	1:G:193:ILE:C	2.23	0.58
1:G:198:PHE:N	1:G:198:PHE:HD1	2.02	0.58
1:I:11:LEU:HD23	1:I:11:LEU:C	2.23	0.58
1:G:183:ALA:O	1:G:187:ALA:CB	2.52	0.58
1:G:146:SER:HB2	1:G:190:ASN:HD22	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:113:ILE:HD13	1:I:116:PRO:HG3	1.86	0.58
2:J:94:SER:HB3	2:J:100:ALA:HA	1.85	0.58
2:J:120:PRO:HB3	2:J:147:PHE:HB3	1.87	0.57
3:D:19:GLU:CB	3:D:75:ARG:NH2	2.67	0.57
1:G:198:PHE:N	1:G:198:PHE:CD1	2.72	0.57
1:I:58:ASP:OD2	1:I:59:GLY:N	2.38	0.57
3:A:141:GLN:O	3:A:145:ARG:HG3	2.04	0.56
1:G:25:TYR:N	1:G:25:TYR:CD2	2.73	0.56
3:D:224:GLN:NE2	3:D:227:ASP:OD2	2.38	0.56
3:A:81:LEU:CD1	5:C:10:LEU:HD11	2.33	0.56
3:A:208:PHE:CD2	3:A:213:ILE:HD12	2.40	0.56
3:A:21:ARG:HG3	3:A:39:ASP:HB2	1.88	0.56
1:G:146:SER:HB2	1:G:190:ASN:ND2	2.20	0.56
1:G:143:THR:CG2	1:G:193:ILE:HG22	2.36	0.56
2:J:216:GLU:OE2	2:J:216:GLU:N	2.38	0.56
1:I:8:SER:O	1:I:104:GLY:O	2.24	0.56
1:I:105:THR:O	1:I:105:THR:HG22	2.05	0.55
1:G:38:GLN:HB2	1:G:44:PRO:CG	2.35	0.55
4:E:29:GLY:HA2	4:E:61:SER:HB2	1.86	0.55
2:J:33:PHE:CE2	2:J:92:ALA:HB3	2.41	0.55
3:A:86:ASN:O	3:A:86:ASN:ND2	2.39	0.55
3:A:13:SER:O	3:A:92:SER:HA	2.07	0.55
2:H:56:ASP:OD1	2:H:57:LYS:N	2.36	0.55
3:A:208:PHE:CD1	3:A:208:PHE:N	2.72	0.55
1:G:56:LYS:O	1:G:62:THR:HG22	2.07	0.55
2:H:39:LEU:HD12	2:H:39:LEU:O	2.06	0.55
3:D:21:ARG:HG3	3:D:39:ASP:HB2	1.88	0.54
1:G:175:ALA:CB	1:G:188:PHE:HE2	2.20	0.54
2:H:168:CYS:SG	2:H:169:THR:N	2.79	0.54
4:B:45:ARG:NH2	4:B:47:GLU:OE1	2.41	0.54
1:G:19:ALA:N	1:G:76:ILE:HD11	2.22	0.54
1:G:123:LEU:HD13	1:G:135:LEU:HD23	1.90	0.54
3:A:208:PHE:N	3:A:208:PHE:HD1	2.06	0.54
3:D:231:VAL:O	3:D:243:LYS:NZ	2.37	0.54
2:J:62:ASP:N	2:J:62:ASP:OD2	2.41	0.54
1:G:197:THR:HB	1:G:199:PHE:CZ	2.43	0.54
1:G:99:LEU:O	2:H:35:TYR:OH	2.26	0.54
3:A:81:LEU:HD11	5:C:10:LEU:CD1	2.34	0.53
1:G:33:PHE:CD1	1:G:65:LEU:CD2	2.91	0.53
1:G:44:PRO:HB2	2:H:102:PHE:HZ	1.73	0.53
1:G:153:VAL:HG22	1:G:177:SER:HB2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:65:LEU:HD11	1:G:72:VAL:HG23	1.90	0.53
1:G:25:TYR:CE2	1:G:70:GLN:HB2	2.39	0.53
2:J:46:ILE:CG2	2:J:47:TYR:H	2.21	0.53
3:A:208:PHE:CE1	3:A:241:PHE:C	2.82	0.53
1:I:52:SER:OG	1:I:55:ASP:OD2	2.27	0.53
1:G:21:LEU:HD21	1:G:74:LEU:CD2	2.38	0.52
1:G:133:VAL:CG2	2:H:125:PHE:HD1	2.22	0.52
2:H:33:PHE:CD1	2:H:48:TYR:HB2	2.45	0.52
1:G:75:LEU:CD2	1:G:75:LEU:N	2.72	0.52
1:G:77:ARG:HD3	1:G:77:ARG:O	2.10	0.52
2:H:20:THR:HG23	2:H:77:THR:HG22	1.92	0.52
1:I:80:GLN:HG2	1:I:81:PRO:HD2	1.92	0.52
3:D:255:GLN:O	3:D:273:ARG:NH2	2.43	0.52
3:D:20:PRO:HG2	3:D:78:LEU:HD12	1.90	0.52
3:D:141:GLN:O	3:D:145:ARG:HG3	2.10	0.52
2:J:119:PRO:HD3	2:J:227:PRO:HB3	1.91	0.52
1:G:25:TYR:CZ	1:G:70:GLN:CB	2.74	0.51
3:D:97:ARG:HD2	3:D:116:PHE:CE2	2.45	0.51
2:H:18:GLN:HA	2:H:78:LEU:O	2.10	0.51
1:G:37:ARG:O	1:G:45:GLU:N	2.40	0.51
2:J:35:TYR:CE2	2:J:45:LEU:HD12	2.45	0.51
1:G:4:VAL:HG22	1:G:25:TYR:HB3	1.93	0.51
2:H:36:ARG:HB3	2:H:46:ILE:HD11	1.93	0.51
1:I:50:ILE:HG23	1:I:65:LEU:HD22	1.91	0.51
2:H:28:SER:OG	2:H:28:SER:O	2.26	0.50
1:G:76:ILE:O	1:G:76:ILE:HG12	2.11	0.50
3:A:189:VAL:HG13	3:A:274:TRP:HB3	1.92	0.50
1:I:19:ALA:HB3	1:I:76:ILE:HB	1.93	0.50
3:D:19:GLU:CG	3:D:75:ARG:NH2	2.75	0.50
1:G:35:TRP:NE1	1:G:74:LEU:HB3	2.27	0.50
1:I:156:THR:OG1	2:J:190:ARG:NH1	2.44	0.50
1:I:97:TYR:HD1	2:J:96:PRO:O	1.94	0.50
3:A:95:LEU:HD21	5:C:10:LEU:HD21	1.94	0.50
1:G:193:ILE:HB	1:G:194:PRO:HD2	1.94	0.49
1:G:33:PHE:CG	1:G:65:LEU:HD21	2.47	0.49
2:J:19:VAL:CG1	2:J:78:LEU:HD21	2.42	0.49
1:I:46:LEU:HD22	2:J:99:GLU:HB2	1.94	0.49
3:A:27:TYR:HE2	3:A:32:GLN:HB2	1.78	0.49
1:G:98:LYS:HB2	2:H:48:TYR:CE1	2.48	0.49
1:G:125:ASP:HB2	1:G:129:SER:HB2	1.95	0.49
4:E:13:HIS:HB2	4:E:21:ASN:HD21	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:6:GLN:HG3	1:G:23:CYS:HB3	1.94	0.49
2:J:41:HIS:N	2:J:41:HIS:ND1	2.60	0.49
3:D:97:ARG:NH1	3:D:114:ASN:OD1	2.46	0.48
1:G:13:VAL:HG22	1:G:13:VAL:O	2.12	0.48
3:A:28:VAL:HG23	3:A:33:PHE:CD1	2.48	0.48
1:I:103:SER:O	2:J:42:GLY:CA	2.62	0.48
1:I:98:LYS:NZ	1:I:98:LYS:CB	2.73	0.48
2:J:48:TYR:CE2	2:J:56:ASP:HB2	2.48	0.48
2:J:92:ALA:HB2	2:J:102:PHE:CD1	2.48	0.48
1:G:65:LEU:CD1	1:G:72:VAL:HG23	2.45	0.47
1:I:192:ILE:CG2	1:I:192:ILE:O	2.55	0.47
2:J:28:SER:OG	2:J:28:SER:O	2.24	0.47
3:D:154:GLU:OE1	3:D:157:ARG:NH1	2.48	0.47
3:D:219:ARG:HH11	3:D:256:ARG:HH12	1.61	0.47
2:H:33:PHE:HD1	2:H:48:TYR:HB2	1.80	0.47
1:G:146:SER:CB	1:G:190:ASN:ND2	2.78	0.47
1:I:124:ARG:HH12	2:J:129:GLU:CD	2.18	0.47
3:A:27:TYR:CE2	3:A:32:GLN:HB2	2.50	0.47
1:G:123:LEU:HD11	1:G:135:LEU:HD23	1.96	0.47
1:I:4:VAL:O	1:I:4:VAL:HG23	2.15	0.47
1:G:75:LEU:N	1:G:75:LEU:HD22	2.31	0.46
2:H:93:SER:N	2:H:101:PHE:O	2.44	0.46
1:G:50:ILE:CG2	1:G:65:LEU:HD13	2.45	0.46
1:I:35:TRP:CZ2	1:I:89:CYS:HB2	2.50	0.46
2:H:17:ARG:HA	2:H:17:ARG:NE	2.31	0.46
2:J:222:GLN:HB3	2:J:224:ARG:HG2	1.98	0.46
2:H:212:TYR:HA	2:H:229:THR:HG22	1.96	0.46
2:J:170:ASP:OD2	2:J:188:SER:OG	2.26	0.46
3:A:138:LYS:HA	3:A:141:GLN:HG3	1.98	0.46
3:D:219:ARG:HD2	3:D:224:GLN:NE2	2.30	0.46
1:I:19:ALA:O	1:I:75:LEU:HD12	2.16	0.46
3:A:37:ASP:OD2	3:A:40:ALA:HB2	2.16	0.45
3:A:61:ASP:O	3:A:65:GLN:HG2	2.15	0.45
1:G:110:ARG:HB3	1:G:141:SER:HB2	1.97	0.45
1:G:193:ILE:C	1:G:193:ILE:CD1	2.85	0.45
2:J:98:THR:HG22	2:J:99:GLU:H	1.82	0.45
1:G:18:ILE:HG23	1:G:18:ILE:O	2.17	0.45
1:G:146:SER:CB	1:G:190:ASN:HD22	2.29	0.45
1:I:166:MET:N	1:I:166:MET:SD	2.89	0.45
2:H:204:HIS:CE1	2:H:235:GLU:HB3	2.52	0.45
1:I:53:ASN:HA	1:I:65:LEU:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:127:PRO:HD3	2:H:140:LEU:HG	1.98	0.45
3:A:185:PRO:HB3	3:A:208:PHE:HB3	1.98	0.45
1:G:57:GLU:HA	1:G:62:THR:CG2	2.43	0.45
1:G:69:SER:HB3	1:G:71:TYR:CD2	2.51	0.45
2:H:43:LEU:HD12	2:H:44:ARG:N	2.31	0.45
2:J:45:LEU:HD23	2:J:58:GLY:HA3	1.99	0.45
1:I:123:LEU:HD11	1:I:135:LEU:HB2	1.99	0.45
1:G:133:VAL:CG2	2:H:125:PHE:CD1	3.00	0.45
1:I:19:ALA:O	1:I:75:LEU:HA	2.17	0.45
1:G:133:VAL:HG12	1:G:176:TRP:CA	2.43	0.45
1:I:148:SER:N	1:I:155:ILE:HD11	2.32	0.45
3:D:219:ARG:HD2	3:D:224:GLN:HE22	1.82	0.44
1:G:66:ASN:HD22	1:G:66:ASN:C	2.19	0.44
1:G:192:ILE:CD1	1:G:192:ILE:N	2.72	0.44
2:H:115:ARG:H	2:H:115:ARG:HG2	1.63	0.44
3:A:107:GLY:O	3:A:169:ARG:NH1	2.50	0.44
2:H:216:GLU:O	2:H:226:LYS:NZ	2.50	0.44
3:A:37:ASP:OD1	3:A:40:ALA:N	2.50	0.44
1:G:133:VAL:HG21	2:H:125:PHE:CD1	2.49	0.44
2:H:34:TRP:O	2:H:45:LEU:HD12	2.16	0.44
3:D:9:TYR:HE1	3:D:99:TYR:HE2	1.66	0.44
2:H:97:GLY:N	5:F:5:VAL:O	2.51	0.44
2:J:212:TYR:HA	2:J:229:THR:HG23	1.99	0.44
2:J:46:ILE:CG2	2:J:47:TYR:N	2.79	0.44
2:H:95:ASP:HB2	2:H:98:THR:OG1	2.18	0.44
1:I:185:ALA:HA	1:I:199:PHE:HE1	1.83	0.44
3:A:131:ARG:HG2	3:A:157:ARG:NH2	2.33	0.44
3:A:96:GLN:OE1	4:B:60:TRP:HB3	2.17	0.44
1:G:19:ALA:H	1:G:76:ILE:HD11	1.83	0.44
3:A:28:VAL:HG23	3:A:33:PHE:HD1	1.83	0.44
2:H:170:ASP:HB3	2:H:187:LEU:HD12	2.00	0.44
3:A:186:LYS:N	3:A:186:LYS:HD2	2.33	0.43
1:G:183:ALA:HB3	1:G:186:ASN:CB	2.33	0.43
1:G:76:ILE:CD1	1:G:76:ILE:N	2.72	0.43
3:A:164:CYS:SG	3:A:165:VAL:N	2.91	0.43
3:A:103:LEU:HD23	3:A:103:LEU:HA	1.90	0.43
3:D:35:GLN:NE2	4:E:53:ASP:OD1	2.51	0.43
4:B:56:PHE:HB3	4:B:62:PHE:CD1	2.53	0.43
1:G:58:ASP:N	1:G:61:PHE:O	2.50	0.43
2:J:70:LYS:HB2	2:J:73:ASN:OD1	2.19	0.43
3:A:156:ARG:O	3:A:160:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:82:THR:N	2:H:85:GLN:OE1	2.52	0.43
1:I:98:LYS:NZ	1:I:98:LYS:HB3	2.33	0.43
4:E:13:HIS:HB2	4:E:21:ASN:ND2	2.34	0.43
1:I:154:TYR:HD2	1:I:176:TRP:NE1	2.17	0.43
3:D:28:VAL:HG23	3:D:33:PHE:CE1	2.54	0.43
4:E:5:PRO:HB3	4:E:30:PHE:HB3	2.00	0.43
1:G:19:ALA:H	1:G:76:ILE:CD1	2.32	0.43
1:G:72:VAL:O	1:G:72:VAL:HG13	2.18	0.43
3:D:19:GLU:HB2	3:D:75:ARG:HH21	1.82	0.42
1:G:175:ALA:HB2	1:G:188:PHE:HE2	1.83	0.42
3:A:9:TYR:HB2	3:A:97:ARG:HB3	2.01	0.42
1:I:35:TRP:CD2	1:I:74:LEU:HD12	2.54	0.42
1:I:97:TYR:HB2	2:J:31:TYR:CZ	2.54	0.42
3:A:188:HIS:O	3:A:204:TRP:HB2	2.19	0.42
4:B:12:ARG:CZ	4:B:22:PHE:HD1	2.32	0.42
3:A:13:SER:HA	3:A:20:PRO:HB3	2.01	0.42
2:J:92:ALA:HB2	2:J:102:PHE:CE1	2.55	0.42
3:D:103:LEU:HA	3:D:103:LEU:HD23	1.87	0.42
4:E:19:LYS:HA	4:E:19:LYS:HD3	1.81	0.42
1:G:123:LEU:HD11	1:G:135:LEU:CD2	2.49	0.42
1:G:144:ASN:OD1	1:G:144:ASN:N	2.52	0.42
2:J:156:TRP:HD1	2:J:167:VAL:HG13	1.85	0.42
1:I:97:TYR:CD1	2:J:96:PRO:C	2.93	0.42
1:I:140:ASP:HB2	1:I:143:THR:HG23	2.01	0.42
3:D:66:LYS:HD3	5:F:2:ALA:HB3	2.02	0.42
3:D:185:PRO:HB3	3:D:208:PHE:HB3	2.01	0.42
3:D:9:TYR:HB2	3:D:97:ARG:HB3	2.01	0.42
1:I:155:ILE:HG22	1:I:175:ALA:HA	2.01	0.42
1:I:92:ALA:HA	1:I:99:LEU:HA	2.02	0.42
3:A:14:ARG:HG3	3:A:14:ARG:O	2.20	0.42
3:A:208:PHE:CD1	3:A:241:PHE:O	2.73	0.42
3:D:47:PRO:HB3	3:D:60:TRP:CZ2	2.55	0.42
2:H:129:GLU:OE1	2:H:132:ILE:HD11	2.20	0.42
1:G:66:ASN:ND2	1:G:66:ASN:C	2.73	0.41
1:I:135:LEU:HD23	1:I:174:VAL:HG22	2.01	0.41
3:A:218:GLN:HB3	3:A:223:ASP:OD1	2.20	0.41
4:B:23:LEU:O	4:B:67:TYR:HA	2.19	0.41
1:G:83:ASP:N	1:G:83:ASP:OD1	2.53	0.41
1:G:95:SER:OG	3:D:62:ARG:NH1	2.53	0.41
1:I:25:TYR:CE1	1:I:70:GLN:HA	2.55	0.41
3:D:22:PHE:HB3	3:D:38:SER:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:21:LEU:HD21	1:I:107:LEU:HD12	2.02	0.41
3:A:58:GLU:OE1	3:A:58:GLU:N	2.44	0.41
3:D:58:GLU:OE1	3:D:58:GLU:N	2.43	0.41
1:G:38:GLN:HA	1:G:44:PRO:HA	2.02	0.41
2:H:35:TYR:CG	2:H:43:LEU:HD11	2.54	0.41
2:H:92:ALA:HB2	2:H:102:PHE:HB3	2.01	0.41
3:D:177:LYS:HB3	3:D:177:LYS:HE2	1.88	0.41
2:J:223:ASP:OD1	2:J:224:ARG:N	2.54	0.41
1:G:36:TYR:OH	2:H:100:ALA:N	2.51	0.41
1:I:90:ALA:HB2	1:I:101:PHE:CE1	2.55	0.41
3:D:13:SER:HA	3:D:20:PRO:HB3	2.01	0.41
4:E:25:CYS:HB2	4:E:39:LEU:HD21	2.02	0.41
4:E:54:LEU:HD11	4:E:62:PHE:CD2	2.39	0.41
2:H:31:TYR:HB2	2:H:94:SER:O	2.20	0.41
2:J:156:TRP:CD1	2:J:167:VAL:HG13	2.55	0.41
4:E:23:LEU:HB2	4:E:70:PHE:CE1	2.55	0.41
1:G:37:ARG:CB	1:G:47:ILE:HD11	2.51	0.41
1:I:113:ILE:HD12	1:I:113:ILE:O	2.21	0.41
3:A:81:LEU:HA	3:A:81:LEU:HD23	1.85	0.40
4:B:23:LEU:HB3	4:B:68:THR:HG22	2.02	0.40
4:E:64:LEU:HA	4:E:64:LEU:HD23	1.89	0.40
4:E:23:LEU:O	4:E:67:TYR:HA	2.21	0.40
1:G:66:ASN:ND2	1:G:66:ASN:O	2.50	0.40
2:H:39:LEU:CG	2:H:39:LEU:O	2.70	0.40
1:I:169:LYS:HE2	1:I:169:LYS:HB3	1.87	0.40
3:D:178:THR:O	3:D:181:ARG:NH1	2.54	0.40
3:A:270:LEU:HA	3:A:270:LEU:HD23	1.89	0.40
2:H:210:GLN:HG2	2:H:212:TYR:CZ	2.57	0.40
4:E:24:ASN:HB3	4:E:65:LEU:HD11	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	197/202 (98%)	185 (94%)	12 (6%)	0	100	100
1	I	189/202 (94%)	179 (95%)	10 (5%)	0	100	100
2	H	236/241 (98%)	227 (96%)	9 (4%)	0	100	100
2	J	236/241 (98%)	228 (97%)	8 (3%)	0	100	100
3	A	271/274 (99%)	264 (97%)	7 (3%)	0	100	100
3	D	271/274 (99%)	263 (97%)	8 (3%)	0	100	100
4	B	96/99 (97%)	94 (98%)	2 (2%)	0	100	100
4	E	95/99 (96%)	93 (98%)	2 (2%)	0	100	100
5	C	8/10 (80%)	8 (100%)	0	0	100	100
5	F	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
All	All	1607/1652 (97%)	1548 (96%)	59 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	G	177/180 (98%)	153 (86%)	24 (14%)	3	20
1	I	171/180 (95%)	154 (90%)	17 (10%)	8	33
2	H	207/209 (99%)	199 (96%)	8 (4%)	32	64
2	J	207/209 (99%)	196 (95%)	11 (5%)	22	55
3	A	229/230 (100%)	217 (95%)	12 (5%)	23	56
3	D	229/230 (100%)	223 (97%)	6 (3%)	46	74
4	B	93/94 (99%)	92 (99%)	1 (1%)	73	88
4	E	92/94 (98%)	92 (100%)	0	100	100
5	C	5/5 (100%)	5 (100%)	0	100	100
5	F	5/5 (100%)	4 (80%)	1 (20%)	1	7
All	All	1415/1436 (98%)	1335 (94%)	80 (6%)	20	53

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	6	GLN
1	G	13	VAL
1	G	18	ILE
1	G	25	TYR
1	G	34	PHE
1	G	55	ASP
1	G	60	ARG
1	G	62	THR
1	G	66	ASN
1	G	70	GLN
1	G	74	LEU
1	G	75	LEU
1	G	76	ILE
1	G	77	ARG
1	G	78	ASP
1	G	127	LYS
1	G	128	SER
1	G	129	SER
1	G	133	VAL
1	G	157	ASP
1	G	190	ASN
1	G	192	ILE
1	G	197	THR
1	G	198	PHE
2	H	18	GLN
2	H	33	PHE
2	H	43	LEU
2	H	102	PHE
2	H	115	ARG
2	H	168	CYS
2	H	181	ASN
2	H	190	ARG
1	I	5	GLU
1	I	13	VAL
1	I	15	GLU
1	I	34	PHE
1	I	37	ARG
1	I	38	GLN
1	I	51	TYR
1	I	74	LEU
1	I	77	ARG
1	I	95	SER

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Mol	Chain	Res	Type
1	I	96	SER
1	I	98	LYS
1	I	105	THR
1	I	115	ASN
1	I	157	ASP
1	I	166	MET
1	I	182	PHE
2	J	18	GLN
2	J	33	PHE
2	J	38	ASP
2	J	41	HIS
2	J	45	LEU
2	J	48	TYR
2	J	87	SER
2	J	184	ARG
2	J	190	ARG
2	J	203	ASN
2	J	239	ARG
3	A	6	ARG
3	A	37	ASP
3	A	119	ASP
3	A	156	ARG
3	A	164	CYS
3	A	180	GLN
3	A	181	ARG
3	A	195	SER
3	A	208	PHE
3	A	212	GLU
3	A	219	ARG
3	A	274	TRP
4	B	47	GLU
3	D	19	GLU
3	D	146	LYS
3	D	152	GLU
3	D	164	CYS
3	D	181	ARG
3	D	196	ASP
5	F	10	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	64	GLN
1	G	70	GLN
1	G	80	GLN
3	D	35	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 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	G	199/202 (98%)	0.33	8 (4%) 38 33	31, 169, 253, 303	0
1	I	193/202 (95%)	0.68	17 (8%) 10 10	97, 168, 225, 308	0
2	H	238/241 (98%)	0.11	11 (4%) 32 28	82, 110, 153, 181	0
2	J	238/241 (98%)	0.02	3 (1%) 77 71	89, 120, 165, 188	0
3	A	273/274 (99%)	-0.08	7 (2%) 56 49	75, 112, 158, 277	0
3	D	273/274 (99%)	-0.17	2 (0%) 87 83	80, 116, 169, 202	0
4	B	98/99 (98%)	-0.29	1 (1%) 82 77	72, 100, 129, 147	0
4	E	97/99 (97%)	-0.13	1 (1%) 82 77	82, 116, 146, 162	0
5	C	10/10 (100%)	0.41	0 100 100	95, 102, 120, 122	0
5	F	10/10 (100%)	-0.16	0 100 100	92, 106, 126, 130	0
All	All	1629/1652 (98%)	0.07	50 (3%) 49 43	31, 121, 197, 308	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	136	PHE	5.2
2	H	79	GLU	5.2
1	G	182	PHE	4.5
2	H	138	ALA	4.4
2	H	137	LYS	4.2
1	I	162	ASP	4.2
1	I	169	LYS	4.1
2	J	139	THR	3.7
3	A	106	ASP	3.5
1	I	197	THR	3.2
3	D	13	SER	3.2
1	I	161	LEU	3.2
1	I	188	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
3	A	220	ASP	3.1
1	I	184	CYS	3.1
1	I	79	SER	2.9
1	I	119	ALA	2.9
3	A	91	GLY	2.9
3	A	92	SER	2.9
1	I	112	ASP	2.8
1	G	200	PRO	2.8
2	H	136	GLN	2.8
2	H	139	THR	2.8
1	G	136	PHE	2.7
1	I	20	SER	2.5
1	G	57	GLU	2.4
2	H	14	GLU	2.4
1	I	123	LEU	2.4
2	H	219	GLU	2.4
1	I	51	TYR	2.4
2	J	239	ARG	2.4
1	I	19	ALA	2.3
3	A	9	TYR	2.3
4	E	28	SER	2.3
2	J	77	THR	2.3
3	D	130	LEU	2.3
2	H	135	THR	2.3
1	I	176	TRP	2.3
1	I	186	ASN	2.2
1	G	188	PHE	2.2
3	A	74	ASP	2.2
2	H	78	LEU	2.2
1	G	77	ARG	2.2
3	A	187	THR	2.2
2	H	192	ARG	2.2
4	B	72	PRO	2.1
1	G	173	ALA	2.1
2	H	57	LYS	2.1
1	I	194	PRO	2.0
1	G	174	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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.