



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:27 PM GMT

PDB ID : 4EUP
Title : The complex between TCR JKF6 and human Class I MHC HLA-A2 presenting the MART-1(27-35)(A27L) peptide
Authors : Hossain, M.; Baker, B.M.
Deposited on : 2012-04-25
Resolution : 2.88 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

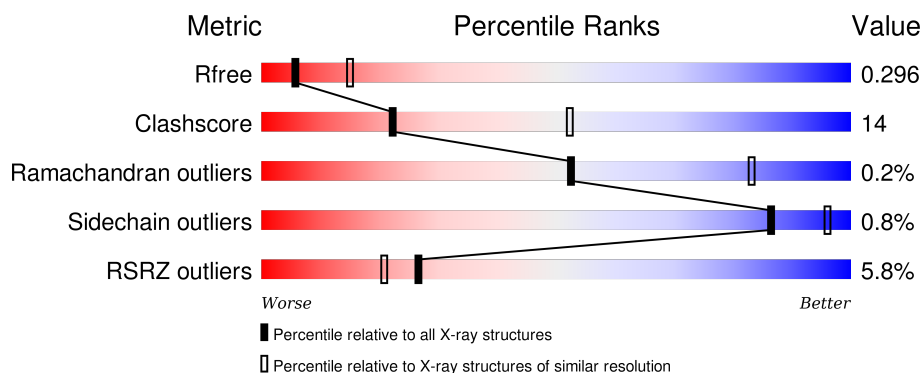
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1945 (2.90-2.86)
Clashscore	102246	2202 (2.90-2.86)
Ramachandran outliers	100387	2149 (2.90-2.86)
Sidechain outliers	100360	2152 (2.90-2.86)
RSRZ outliers	91569	1950 (2.90-2.86)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	275	<div> <div>3%</div> <div>73%</div> <div>23%</div> <div>• •</div> </div>
1	D	275	<div> <div>4%</div> <div>72%</div> <div>23%</div> <div>• •</div> </div>
2	B	100	<div> <div>2%</div> <div>68%</div> <div>32%</div> </div>
2	E	100	<div> <div>%</div> <div>83%</div> <div>17%</div> </div>
3	C	9	<div> <div>56%</div> <div>44%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	9	<div><div></div><div>78%</div><div>22%</div></div>
4	G	206	<div><div>11%</div><div></div><div>77%</div><div>20%</div><div>••</div></div>
4	I	206	<div><div>15%</div><div></div><div>76%</div><div>21%</div><div>•</div></div>
5	H	243	<div><div>4%</div><div></div><div>76%</div><div>23%</div><div>•</div></div>
5	J	243	<div><div>4%</div><div></div><div>71%</div><div>28%</div><div>•</div></div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13167 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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	265	Total	C	N	O	S	0	0	0
			2164	1359	394	402	9			
1	A	267	Total	C	N	O	S	0	0	0
			2183	1369	399	406	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	100	Total	C	N	O	S	0	0	0
			836	533	141	158	4			
2	B	100	Total	C	N	O	S	0	0	0
			836	533	141	158	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	0	MET	-	INITIATING METHIONINE	UNP P61769
B	0	MET	-	INITIATING METHIONINE	UNP P61769

- Molecule 3 is a protein called Melanoma antigen recognized by T-cells 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	9	Total	C	N	O	0	0	0
			60	40	9	11			
3	C	9	Total	C	N	O	0	0	0
			60	40	9	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	2	LEU	ALA	ENGINEERED MUTATION	UNP Q16655
C	2	LEU	ALA	ENGINEERED MUTATION	UNP Q16655

- Molecule 4 is a protein called JKF6 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	201	Total	C	N	O	S	0	0	0
			1537	955	255	319	8			
4	I	201	Total	C	N	O	S	0	0	0
			1537	955	255	319	8			

- Molecule 5 is a protein called JKF6 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	H	241	Total	C	N	O	S	0	0	0
			1938	1229	330	370	9			
5	J	240	Total	C	N	O	S	0	0	0
			1930	1225	329	367	9			

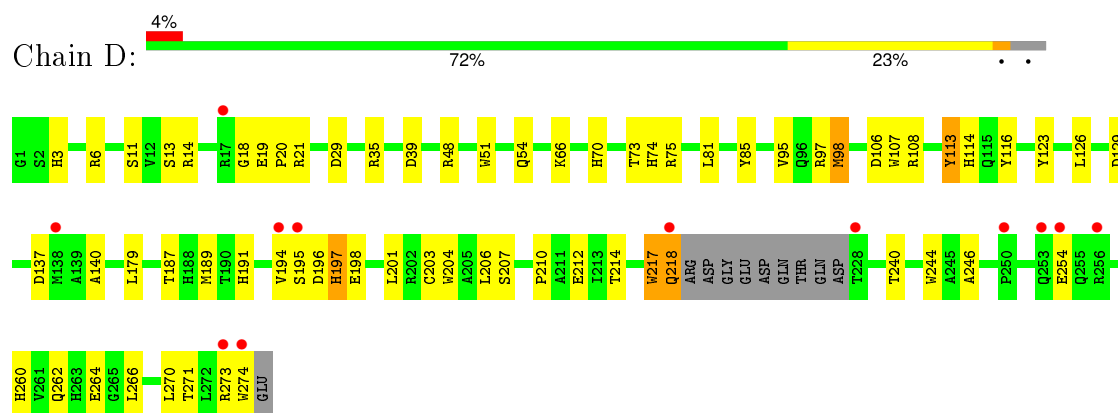
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	13	Total	O	0	0
			13	13		
6	E	9	Total	O	0	0
			9	9		
6	A	10	Total	O	0	0
			10	10		
6	B	12	Total	O	0	0
			12	12		
6	G	10	Total	O	0	0
			10	10		
6	I	5	Total	O	0	0
			5	5		
6	H	13	Total	O	0	0
			13	13		
6	J	14	Total	O	0	0
			14	14		

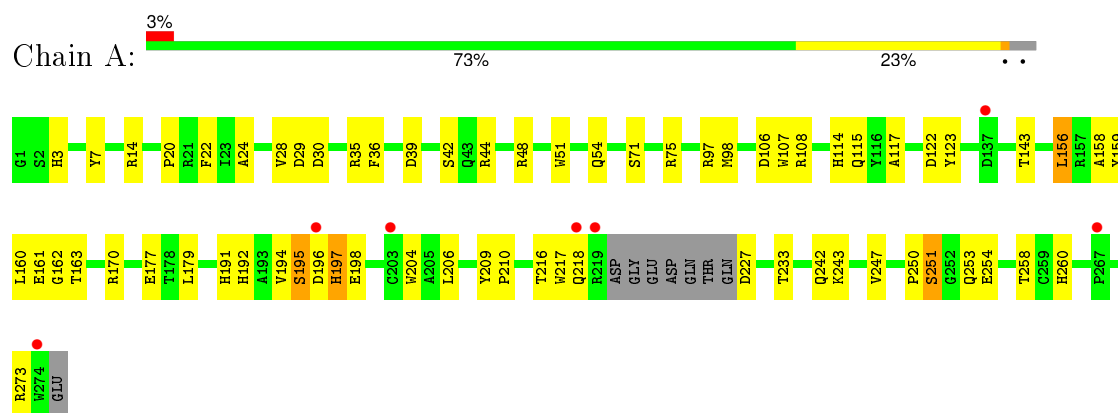
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

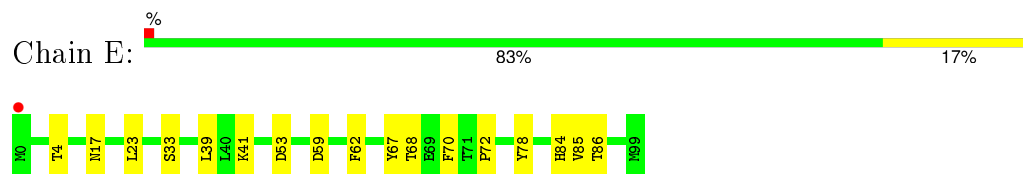
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



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- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin





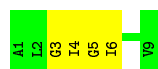
- Molecule 3: Melanoma antigen recognized by T-cells 1

Chain F: 78% 22%



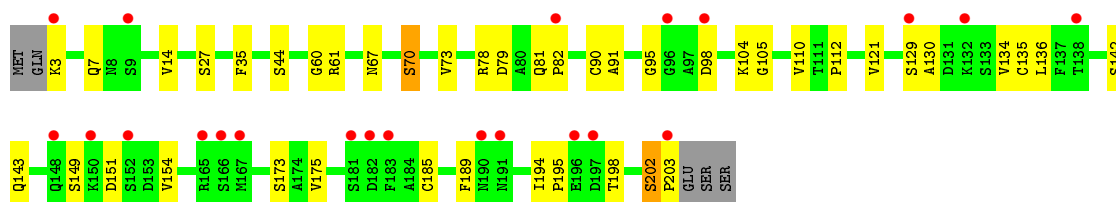
- Molecule 3: Melanoma antigen recognized by T-cells 1

Chain C: 56% 44%



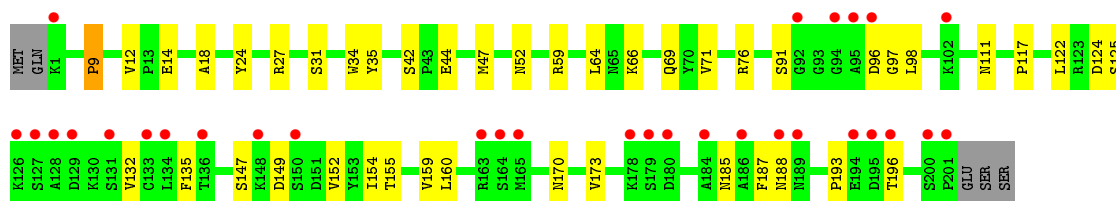
- Molecule 4: JFK6 alpha chain

Chain G: 11% 77% 20%



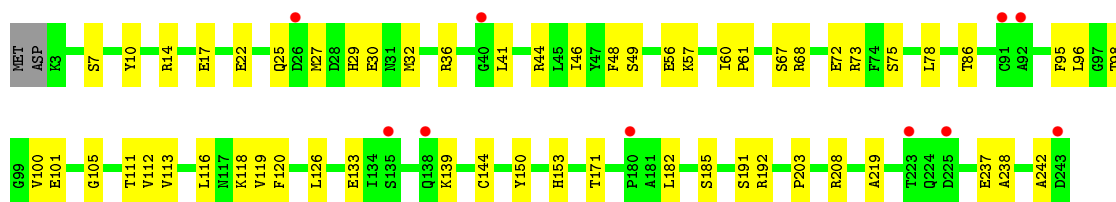
- Molecule 4: JFK6 alpha chain

Chain I: 15% 76% 21%

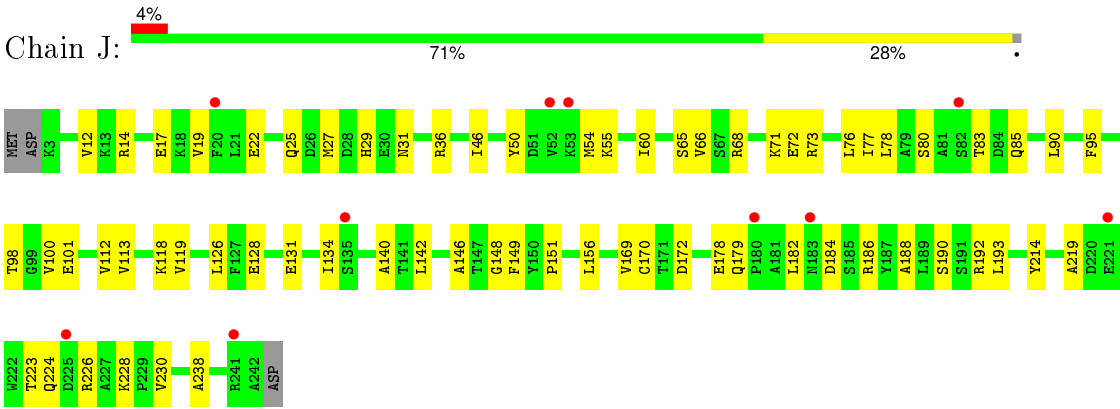


- Molecule 5: JFK6 beta chain

Chain H: 4% 76% 23%



- Molecule 5: JFK6 beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.66 Å 53.70 Å 150.51 Å 90.00° 112.46° 90.00°	Depositor
Resolution (Å)	20.00 – 2.88 19.94 – 2.88	Depositor EDS
% Data completeness (in resolution range)	92.6 (20.00-2.88) 92.6 (19.94-2.88)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.88 Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.235 , 0.295 0.242 , 0.296	Depositor DCC
R_{free} test set	2170 reflections (5.71%)	DCC
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.089	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 59.3	EDS
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 39995 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13167	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.21 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.1727e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [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	A	0.79	0/2247	0.80	2/3049 (0.1%)
1	D	0.78	2/2228 (0.1%)	0.90	7/3024 (0.2%)
2	B	0.84	1/859 (0.1%)	0.78	0/1162
2	E	0.81	0/859	0.76	0/1162
3	C	0.78	0/59	0.93	0/78
3	F	0.81	0/59	1.05	0/78
4	G	0.69	2/1569 (0.1%)	0.78	3/2128 (0.1%)
4	I	0.65	0/1569	0.77	3/2128 (0.1%)
5	H	0.61	0/1988	0.68	1/2696 (0.0%)
5	J	0.66	0/1980	0.72	1/2685 (0.0%)
All	All	0.72	5/13417 (0.0%)	0.78	17/18190 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	90	CYS	CB-SG	-7.34	1.69	1.82
4	G	202	SER	CB-OG	6.79	1.51	1.42
1	D	212	GLU	CG-CD	5.47	1.60	1.51
1	D	217	TRP	C-N	-5.29	1.21	1.34
2	B	77	GLU	CB-CG	5.24	1.62	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	217	TRP	O-C-N	-13.31	101.41	122.70
1	D	217	TRP	C-N-CA	9.67	145.89	121.70
1	D	217	TRP	CA-C-N	9.51	138.13	117.20
4	I	9	PRO	CA-N-CD	-8.48	99.63	111.50
4	I	97	GLY	N-CA-C	-8.04	93.00	113.10
4	G	135	CYS	CA-CB-SG	-6.67	102.00	114.00
4	G	185	CYS	CA-CB-SG	-6.54	102.23	114.00
1	D	98	MET	CB-CG-SD	-6.31	93.48	112.40
1	D	197	HIS	N-CA-C	-6.23	94.17	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	129	ASP	CB-CA-C	6.15	122.69	110.40
1	A	273	ARG	N-CA-C	-6.11	94.51	111.00
1	A	122	ASP	CB-CG-OD2	5.87	123.58	118.30
1	D	126	LEU	CB-CA-C	5.73	121.09	110.20
5	J	214	TYR	CB-CA-C	-5.48	99.44	110.40
5	H	144	CYS	CA-CB-SG	-5.19	104.65	114.00
4	I	98	LEU	N-CA-CB	5.19	120.78	110.40
4	G	90	CYS	CA-CB-SG	-5.02	104.96	114.00

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	2183	0	2049	61	0
1	D	2164	0	2031	72	0
2	B	836	0	803	32	0
2	E	836	0	803	19	0
3	C	60	0	73	6	0
3	F	60	0	73	4	0
4	G	1537	0	1448	41	0
4	I	1537	0	1451	37	0
5	H	1938	0	1862	59	0
5	J	1930	0	1858	59	0
6	A	10	0	0	0	0
6	B	12	0	0	0	0
6	D	13	0	0	1	0
6	E	9	0	0	0	0
6	G	10	0	0	0	0
6	H	13	0	0	0	0
6	I	5	0	0	0	0
6	J	14	0	0	1	0
All	All	13167	0	12451	352	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 14.

All (352) 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:D:196:ASP:HB3	1:D:197:HIS:CD2	1.39	1.55
1:D:196:ASP:CB	1:D:197:HIS:HD2	1.47	1.26
4:G:7:GLN:NE2	4:G:105:GLY:H	1.36	1.24
1:D:196:ASP:CB	1:D:197:HIS:CD2	2.21	1.21
1:D:35:ARG:HD3	2:E:53:ASP:OD1	1.42	1.17
5:H:86:THR:HG22	5:H:112:VAL:H	1.07	1.13
4:G:202:SER:HB2	4:G:203:PRO:HD2	1.33	1.07
4:G:7:GLN:NE2	4:G:105:GLY:N	2.04	1.05
2:B:17:ASN:ND2	2:B:72:PRO:HB2	1.74	1.03
4:I:149:ASP:HB3	4:I:152:VAL:HG12	1.46	0.97
1:D:191:HIS:HB3	1:D:274:TRP:CH2	1.99	0.97
5:H:86:THR:CG2	5:H:112:VAL:H	1.80	0.92
2:B:17:ASN:HD22	2:B:72:PRO:HB2	1.32	0.92
5:H:86:THR:HG22	5:H:112:VAL:N	1.85	0.92
2:B:22:PHE:CE2	2:B:69:GLU:HG2	2.04	0.92
1:A:3:HIS:HD2	1:A:29:ASP:OD2	1.55	0.90
4:G:7:GLN:HE21	4:G:105:GLY:N	1.69	0.89
2:E:23:LEU:HB2	2:E:70:PHE:CD2	2.08	0.89
1:A:191:HIS:NE2	1:A:254:GLU:OE2	2.07	0.87
4:I:52:ASN:HD21	4:I:66:LYS:H	1.23	0.87
1:A:194:VAL:O	1:A:195:SER:HB3	1.76	0.86
1:D:254:GLU:HB3	1:D:274:TRP:HD1	1.41	0.85
1:D:254:GLU:HB3	1:D:274:TRP:CD1	2.12	0.83
4:G:7:GLN:NE2	4:G:105:GLY:CA	2.41	0.83
1:D:189:MET:SD	1:D:201:LEU:HD23	2.18	0.83
1:D:214:THR:HG23	1:D:262:GLN:HB2	1.60	0.81
1:A:197:HIS:HA	1:A:251:SER:HB2	1.62	0.80
1:D:18:GLY:O	1:D:19:GLU:C	2.19	0.80
5:J:182:LEU:HD23	5:J:184:ASP:H	1.44	0.80
1:D:196:ASP:CG	1:D:197:HIS:HD2	1.84	0.79
4:G:7:GLN:HE21	4:G:105:GLY:CA	1.96	0.79
1:D:191:HIS:CB	1:D:274:TRP:CH2	2.64	0.79
5:H:49:SER:OG	5:H:68:ARG:HD3	1.83	0.78
5:J:182:LEU:HD21	5:J:184:ASP:OD1	1.84	0.77
5:H:7:SER:HB3	5:H:22:GLU:HB2	1.65	0.77
4:I:14:GLU:HB3	4:I:111:ASN:HD21	1.50	0.76
1:D:196:ASP:HB3	1:D:197:HIS:CG	2.16	0.76
1:D:3:HIS:HD2	1:D:29:ASP:OD2	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:149:ASP:HB3	4:I:152:VAL:CG1	2.17	0.75
4:I:185:ASN:HB2	4:I:188:ASN:HD22	1.52	0.75
4:G:189:PHE:HB3	4:G:194:ILE:HD11	1.69	0.74
1:D:187:THR:HG21	1:D:270:LEU:HD23	1.69	0.74
1:D:35:ARG:HD3	2:E:53:ASP:CG	2.08	0.74
5:J:36:ARG:HH12	5:J:85:GLN:HA	1.54	0.73
4:G:67:ASN:O	4:G:70:SER:O	2.05	0.73
4:G:202:SER:HB2	4:G:203:PRO:CD	2.16	0.72
4:I:193:PRO:O	4:I:196:THR:HG22	1.89	0.72
4:I:96:ASP:OD1	5:J:100:VAL:HG22	1.89	0.72
1:A:48:ARG:NH2	2:B:53:ASP:OD1	2.22	0.72
2:B:3:ARG:HH21	2:B:61:SER:HB3	1.54	0.72
4:G:7:GLN:HE22	4:G:105:GLY:H	1.34	0.71
4:G:149:SER:HB2	4:G:154:VAL:HG13	1.70	0.71
2:B:33:SER:HB3	2:B:62:PHE:CE2	2.25	0.71
4:G:202:SER:CB	4:G:203:PRO:HD2	2.16	0.70
4:G:195:PRO:O	4:G:198:THR:HG22	1.90	0.70
4:I:147:SER:HB3	4:I:154:ILE:HD13	1.73	0.70
1:D:35:ARG:HE	1:D:48:ARG:HD3	1.57	0.69
5:H:25:GLN:NE2	5:H:29:HIS:H	1.89	0.69
1:D:35:ARG:CD	2:E:53:ASP:OD1	2.30	0.69
1:D:35:ARG:CD	2:E:53:ASP:CG	2.61	0.68
4:G:189:PHE:O	4:G:194:ILE:HD12	1.93	0.68
5:J:83:THR:HA	5:J:112:VAL:CG2	2.23	0.68
5:H:171:THR:HG23	5:H:191:SER:HB2	1.75	0.68
1:A:206:LEU:HD22	1:A:242:GLN:HG2	1.75	0.68
1:A:194:VAL:O	1:A:195:SER:CB	2.42	0.67
2:E:4:THR:HG22	2:E:86:THR:OG1	1.94	0.67
5:J:126:LEU:HD22	5:J:238:ALA:HB2	1.74	0.67
4:G:61:ARG:NH1	4:G:79:ASP:O	2.28	0.67
4:I:155:THR:HG21	5:J:192:ARG:HH12	1.59	0.67
5:J:182:LEU:HD23	5:J:184:ASP:N	2.10	0.66
1:D:214:THR:CG2	1:D:262:GLN:HB2	2.25	0.65
5:J:83:THR:HA	5:J:112:VAL:HG23	1.78	0.65
5:H:32:MET:SD	5:H:68:ARG:NH1	2.65	0.65
1:A:197:HIS:HA	1:A:251:SER:CB	2.27	0.65
2:B:51:HIS:HA	2:B:65:LEU:O	1.97	0.65
4:G:189:PHE:O	4:G:194:ILE:CD1	2.45	0.64
4:I:52:ASN:ND2	4:I:66:LYS:H	1.92	0.64
2:B:17:ASN:ND2	2:B:97:ARG:HH22	1.95	0.64
1:D:97:ARG:HH21	1:D:114:HIS:CE1	2.15	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:81:GLN:HB3	4:G:82:PRO:HD2	1.79	0.64
1:A:196:ASP:O	1:A:197:HIS:C	2.35	0.64
5:H:57:LYS:HB3	5:H:61:PRO:HG3	1.80	0.64
2:E:4:THR:HG22	2:E:86:THR:CB	2.28	0.63
1:A:197:HIS:CA	1:A:251:SER:HB2	2.29	0.63
1:D:35:ARG:NE	1:D:48:ARG:HH21	1.96	0.63
4:G:112:PRO:HD2	4:G:142:SER:OG	1.98	0.63
1:A:197:HIS:HB3	1:A:251:SER:HB3	1.81	0.62
5:H:126:LEU:HD11	5:H:238:ALA:HB2	1.79	0.62
2:B:7:ILE:HG12	2:B:82:VAL:HG21	1.81	0.62
5:J:228:LYS:HD3	5:J:230:VAL:HG13	1.82	0.62
4:I:135:PHE:HB2	4:I:187:PHE:CE2	2.35	0.61
4:I:14:GLU:HB3	4:I:111:ASN:ND2	2.15	0.61
1:D:210:PRO:HG3	1:D:264:GLU:OE1	2.00	0.61
4:G:104:LYS:NZ	5:J:80:SER:OG	2.32	0.61
1:A:196:ASP:CG	1:A:196:ASP:O	2.38	0.60
5:J:219:ALA:HB3	6:J:312:HOH:O	2.02	0.60
5:H:44:ARG:HH11	5:H:60:ILE:CD1	2.15	0.60
3:F:7:LEU:HD11	5:H:95:PHE:CB	2.31	0.60
4:G:189:PHE:HB3	4:G:194:ILE:CD1	2.31	0.60
4:I:122:LEU:HB3	5:J:128:GLU:O	2.02	0.60
5:H:60:ILE:O	5:H:60:ILE:HG22	2.01	0.59
5:J:65:SER:HB2	5:J:77:ILE:HB	1.82	0.59
3:C:4:ILE:HA	5:J:98:THR:HB	1.85	0.59
5:J:101:GLU:N	5:J:101:GLU:OE1	2.35	0.59
1:A:206:LEU:CD2	1:A:242:GLN:HG2	2.32	0.59
4:I:9:PRO:O	4:I:9:PRO:HD2	2.01	0.59
2:E:70:PHE:CE1	2:E:72:PRO:HD3	2.37	0.58
5:J:36:ARG:HG3	5:J:46:ILE:HD11	1.86	0.58
1:A:158:ALA:O	1:A:161:GLU:O	2.21	0.58
5:J:12:VAL:HG22	5:J:151:PRO:HG3	1.86	0.58
4:G:151:ASP:HB3	4:G:154:VAL:HG12	1.84	0.58
1:A:162:GLY:O	1:A:163:THR:C	2.41	0.57
1:D:13:SER:HA	1:D:20:PRO:HB3	1.86	0.57
1:D:35:ARG:HE	1:D:48:ARG:HH21	1.52	0.57
1:D:191:HIS:CB	1:D:274:TRP:CZ2	2.86	0.57
1:D:75:ARG:HG2	6:D:303:HOH:O	2.04	0.57
5:J:50:TYR:CZ	5:J:54:MET:HG2	2.40	0.57
5:H:49:SER:OG	5:H:68:ARG:CD	2.52	0.57
1:D:106:ASP:OD2	1:D:108:ARG:N	2.32	0.57
2:B:17:ASN:CG	2:B:97:ARG:HH22	2.08	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:ARG:HB2	1:D:116:TYR:CD2	2.41	0.56
2:B:17:ASN:HD21	2:B:97:ARG:HH22	1.53	0.56
1:D:254:GLU:CB	1:D:274:TRP:CD1	2.88	0.56
1:D:14:ARG:NH2	1:D:39:ASP:OD1	2.37	0.56
4:I:96:ASP:OD2	5:J:31:ASN:ND2	2.39	0.55
2:E:4:THR:HG22	2:E:86:THR:HB	1.87	0.55
5:J:22:GLU:HB3	5:J:73:ARG:NH1	2.21	0.55
1:A:42:SER:O	1:A:44:ARG:HG2	2.06	0.55
1:A:51:TRP:O	1:A:54:GLN:HG2	2.07	0.55
5:H:208:ARG:HG3	5:H:237:GLU:HG2	1.87	0.55
4:G:35:PHE:CE1	5:H:101:GLU:HB3	2.41	0.55
1:D:196:ASP:CG	1:D:197:HIS:CD2	2.71	0.55
1:D:266:LEU:HD13	1:D:270:LEU:HD13	1.89	0.55
1:D:191:HIS:CG	1:D:274:TRP:CZ2	2.96	0.54
1:D:35:ARG:HD2	2:E:53:ASP:CG	2.27	0.54
3:F:7:LEU:HD11	5:H:95:PHE:HB3	1.89	0.54
4:I:159:VAL:HG22	4:I:170:ASN:ND2	2.22	0.54
1:A:51:TRP:CE2	1:A:179:LEU:HD11	2.43	0.54
5:H:126:LEU:C	5:H:126:LEU:HD12	2.27	0.54
1:A:22:PHE:CD2	1:A:71:SER:HB3	2.42	0.54
1:A:29:ASP:O	1:A:30:ASP:HB2	2.08	0.54
1:A:3:HIS:CD2	1:A:29:ASP:OD2	2.47	0.53
5:H:36:ARG:HG3	5:H:46:ILE:HD11	1.91	0.53
5:H:44:ARG:HH11	5:H:60:ILE:HD12	1.73	0.53
1:D:196:ASP:HB3	1:D:197:HIS:HD2	0.87	0.53
4:I:185:ASN:HB2	4:I:188:ASN:ND2	2.23	0.53
1:D:260:HIS:ND1	1:D:271:THR:HG22	2.24	0.53
5:H:22:GLU:HB3	5:H:73:ARG:HH11	1.73	0.52
2:B:27:VAL:O	2:B:63:TYR:HA	2.09	0.52
4:G:81:GLN:HB3	4:G:82:PRO:CD	2.39	0.52
1:D:11:SER:HB3	1:D:95:VAL:HG12	1.90	0.52
1:A:204:TRP:HH2	2:B:99:MET:C	2.12	0.52
1:A:197:HIS:N	1:A:197:HIS:ND1	2.56	0.52
5:H:25:GLN:HG2	5:H:27:MET:H	1.74	0.52
5:J:182:LEU:CD2	5:J:184:ASP:C	2.78	0.52
5:J:146:ALA:O	5:J:149:PHE:HE1	1.93	0.52
1:D:196:ASP:OD2	1:D:197:HIS:CD2	2.63	0.52
1:D:273:ARG:O	1:D:274:TRP:HB3	2.10	0.52
1:D:20:PRO:HD2	1:D:75:ARG:HG3	1.92	0.52
2:E:33:SER:HB3	2:E:62:PHE:CE2	2.45	0.52
5:J:134:ILE:HD11	5:J:140:ALA:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:151:ASP:HB3	4:G:154:VAL:CG1	2.39	0.51
5:J:119:VAL:O	5:J:226:ARG:NH2	2.42	0.51
1:A:97:ARG:CZ	3:C:6:ILE:HG23	2.40	0.51
4:G:173:SER:OG	5:H:192:ARG:HD3	2.10	0.51
1:A:216:THR:OG1	1:A:260:HIS:HB2	2.10	0.51
4:G:175:VAL:HG23	5:H:192:ARG:HE	1.74	0.51
5:H:14:ARG:O	5:H:17:GLU:HB2	2.10	0.51
1:A:51:TRP:CZ2	1:A:179:LEU:HD11	2.45	0.51
4:G:98:ASP:OD2	5:H:100:VAL:HG11	2.11	0.51
1:D:51:TRP:O	1:D:54:GLN:HG2	2.10	0.51
4:G:44:SER:CB	5:H:105:GLY:O	2.58	0.51
2:B:50:GLU:HG3	2:B:51:HIS:H	1.75	0.51
5:J:146:ALA:O	5:J:188:ALA:HA	2.11	0.51
5:J:126:LEU:HD21	5:J:142:LEU:HD23	1.92	0.50
4:G:98:ASP:OD2	5:H:100:VAL:CG1	2.59	0.50
1:A:233:THR:OG1	1:A:243:LYS:HE2	2.11	0.50
5:H:30:GLU:HG3	5:H:96:LEU:HB3	1.92	0.50
5:H:78:LEU:HD12	5:H:78:LEU:N	2.26	0.50
2:B:22:PHE:CD2	2:B:69:GLU:HG2	2.46	0.50
5:H:48:PHE:CZ	5:H:56:GLU:HB2	2.47	0.50
2:B:50:GLU:CG	2:B:51:HIS:H	2.25	0.50
1:A:35:ARG:HD3	1:A:48:ARG:HD3	1.93	0.50
1:A:106:ASP:OD2	1:A:107:TRP:N	2.45	0.50
1:A:24:ALA:HB3	1:A:36:PHE:HB3	1.93	0.50
1:D:95:VAL:HG21	1:D:116:TYR:OH	2.11	0.50
5:H:67:SER:HB3	5:H:75:SER:HB2	1.93	0.50
1:A:163:THR:CG2	4:I:66:LYS:NZ	2.75	0.49
5:J:169:VAL:HG23	5:J:193:LEU:HD13	1.93	0.49
4:I:64:LEU:HD13	4:I:71:VAL:HG22	1.94	0.49
2:B:33:SER:HB3	2:B:62:PHE:CD2	2.47	0.49
5:J:12:VAL:CG2	5:J:151:PRO:HG3	2.41	0.49
2:E:84:HIS:CE1	2:E:86:THR:HG23	2.47	0.49
1:A:250:PRO:HG2	1:A:253:GLN:HG3	1.95	0.49
4:G:7:GLN:HE21	4:G:105:GLY:C	2.15	0.49
5:H:22:GLU:OE1	5:H:73:ARG:HD2	2.11	0.49
5:J:169:VAL:CG2	5:J:193:LEU:HD13	2.42	0.49
5:H:126:LEU:HD11	5:H:238:ALA:CB	2.43	0.49
1:A:106:ASP:OD2	1:A:108:ARG:N	2.33	0.49
5:J:25:GLN:NE2	5:J:29:HIS:H	2.11	0.49
4:G:194:ILE:CG2	4:G:198:THR:HG21	2.43	0.49
5:J:25:GLN:NE2	5:J:27:MET:O	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:PRO:HD2	1:A:75:ARG:HG3	1.93	0.49
2:B:54:LEU:HD11	2:B:62:PHE:CD1	2.48	0.49
5:H:203:PRO:HB3	5:H:242:ALA:HB2	1.95	0.49
4:G:7:GLN:HE22	4:G:105:GLY:N	1.97	0.48
3:F:4:ILE:HA	5:H:98:THR:HB	1.94	0.48
2:B:33:SER:HB3	2:B:62:PHE:CZ	2.48	0.48
1:A:97:ARG:HA	1:A:115:GLN:O	2.14	0.48
5:H:44:ARG:NH1	5:H:60:ILE:HD12	2.28	0.48
1:A:106:ASP:C	1:A:106:ASP:OD2	2.52	0.48
2:B:17:ASN:HD21	2:B:72:PRO:HB2	1.71	0.47
1:D:66:LYS:O	1:D:70:HIS:CD2	2.67	0.47
5:J:131:GLU:HA	5:J:134:ILE:HB	1.95	0.47
2:B:4:THR:HG22	2:B:86:THR:HB	1.95	0.47
1:A:35:ARG:HD2	1:A:48:ARG:HH21	1.80	0.47
1:D:51:TRP:CE2	1:D:179:LEU:HD11	2.49	0.47
5:J:223:THR:O	5:J:224:GLN:HB2	2.14	0.47
1:D:18:GLY:O	1:D:19:GLU:O	2.32	0.47
5:J:156:LEU:HD23	5:J:156:LEU:C	2.35	0.47
5:H:25:GLN:NE2	5:H:29:HIS:N	2.60	0.46
4:I:31:SER:HB2	4:I:91:SER:OG	2.15	0.46
4:I:160:LEU:HB3	5:J:170:CYS:HB2	1.97	0.46
1:D:196:ASP:HB3	1:D:197:HIS:CA	2.46	0.46
1:A:206:LEU:HD22	1:A:242:GLN:CG	2.45	0.46
1:D:73:THR:HG22	1:D:74:HIS:HD2	1.80	0.46
5:J:126:LEU:HD22	5:J:238:ALA:CB	2.44	0.46
5:H:101:GLU:N	5:H:101:GLU:OE1	2.48	0.46
5:H:116:LEU:O	5:H:119:VAL:HG23	2.16	0.46
4:G:129:SER:O	4:G:130:ALA:C	2.54	0.46
1:D:6:ARG:NE	1:D:113:TYR:OH	2.43	0.46
5:J:60:ILE:HG23	5:J:60:ILE:O	2.16	0.46
2:B:50:GLU:HG3	2:B:51:HIS:N	2.30	0.46
4:G:7:GLN:NE2	4:G:105:GLY:HA2	2.29	0.46
5:J:148:GLY:O	5:J:186:ARG:HD3	2.15	0.46
1:A:227:ASP:O	1:A:247:VAL:HA	2.15	0.46
1:A:197:HIS:CA	1:A:251:SER:CB	2.93	0.45
5:H:44:ARG:NH1	5:H:60:ILE:CD1	2.79	0.45
1:A:196:ASP:O	1:A:198:GLU:N	2.50	0.45
4:I:24:TYR:CE2	4:I:69:GLN:HG2	2.51	0.45
5:J:178:GLU:O	5:J:179:GLN:HG3	2.16	0.45
5:J:36:ARG:NH1	5:J:85:GLN:HA	2.28	0.45
4:G:3:LYS:HA	4:G:27:SER:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:J:113:VAL:CG2	5:J:118:LYS:HD2	2.45	0.45
2:B:22:PHE:CE2	2:B:69:GLU:CG	2.91	0.45
5:H:41:LEU:HD13	5:H:44:ARG:CZ	2.47	0.45
4:I:24:TYR:CZ	4:I:69:GLN:HA	2.52	0.45
1:D:194:VAL:O	1:D:195:SER:C	2.55	0.45
5:H:32:MET:CG	5:H:68:ARG:HH12	2.29	0.45
5:J:66:VAL:HG12	5:J:76:LEU:HD12	1.99	0.45
2:E:17:ASN:HA	2:E:72:PRO:O	2.17	0.45
1:A:194:VAL:O	1:A:194:VAL:HG12	2.15	0.45
4:I:147:SER:HB3	4:I:154:ILE:CD1	2.43	0.45
1:D:203:CYS:HB2	1:D:217:TRP:CZ2	2.52	0.45
1:A:192:HIS:CE1	2:B:98:ASP:HB3	2.52	0.45
5:H:29:HIS:HD2	5:H:95:PHE:CE1	2.35	0.45
2:B:70:PHE:CZ	2:B:72:PRO:HG3	2.51	0.44
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.99	0.44
4:G:60:GLY:O	4:G:78:ARG:NH2	2.43	0.44
4:I:35:TYR:HA	4:I:44:GLU:O	2.16	0.44
1:D:244:TRP:HZ3	1:D:246:ALA:HB3	1.81	0.44
2:E:85:VAL:HG12	2:E:85:VAL:O	2.17	0.44
1:D:106:ASP:OD2	1:D:108:ARG:HB2	2.17	0.44
1:A:97:ARG:HH21	1:A:114:HIS:CE1	2.36	0.44
4:I:96:ASP:CG	5:J:100:VAL:HG22	2.38	0.44
1:D:194:VAL:HG22	1:D:198:GLU:HG3	2.00	0.44
5:J:68:ARG:NH2	5:J:71:LYS:O	2.45	0.44
5:J:95:PHE:O	5:J:100:VAL:HG21	2.18	0.44
5:H:118:LYS:O	5:H:120:PHE:HD2	2.01	0.44
1:D:19:GLU:HA	1:D:20:PRO:HD3	1.81	0.43
5:H:22:GLU:HB3	5:H:73:ARG:NH1	2.33	0.43
1:D:207:SER:HA	1:D:240:THR:OG1	2.17	0.43
1:D:81:LEU:O	1:D:85:TYR:HD2	2.02	0.43
3:C:4:ILE:HA	5:J:98:THR:CB	2.47	0.43
2:E:23:LEU:HB2	2:E:70:PHE:CE2	2.52	0.43
5:H:30:GLU:HG2	5:H:96:LEU:H	1.82	0.43
1:A:159:TYR:CD1	3:C:3:GLY:HA3	2.54	0.43
1:A:159:TYR:CE1	3:C:3:GLY:HA3	2.54	0.43
5:H:113:VAL:CG1	5:H:150:TYR:OH	2.67	0.43
1:D:204:TRP:CE3	1:D:206:LEU:HD21	2.53	0.43
4:G:121:VAL:HA	4:G:136:LEU:O	2.19	0.43
5:J:19:VAL:HB	5:J:78:LEU:HB2	2.00	0.43
4:G:189:PHE:O	4:G:194:ILE:HD11	2.18	0.43
1:D:106:ASP:OD2	1:D:107:TRP:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:THR:CG2	1:A:260:HIS:CE1	3.02	0.43
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.54	0.43
5:H:86:THR:HG22	5:H:111:THR:HA	2.01	0.43
1:D:106:ASP:C	1:D:106:ASP:OD2	2.57	0.43
2:B:41:LYS:O	2:B:42:ASN:C	2.57	0.43
1:D:196:ASP:HB3	1:D:197:HIS:HA	1.99	0.43
4:G:134:VAL:HG13	4:G:175:VAL:HG13	2.00	0.43
5:H:113:VAL:HG11	5:H:150:TYR:CE1	2.54	0.43
4:I:12:VAL:HG21	4:I:18:ALA:HB2	2.00	0.43
5:J:55:LYS:HE2	5:J:55:LYS:HB3	1.84	0.43
1:D:254:GLU:CB	1:D:274:TRP:HD1	2.20	0.43
4:I:59:ARG:HA	4:I:76:ARG:NH1	2.34	0.43
1:A:156:LEU:O	1:A:160:LEU:HG	2.19	0.42
1:A:28:VAL:HG12	1:A:29:ASP:HB2	2.01	0.42
4:G:14:VAL:O	4:G:110:VAL:HA	2.20	0.42
4:I:34:TRP:HB2	4:I:47:MET:HB3	2.01	0.42
1:D:51:TRP:CZ2	1:D:179:LEU:HD11	2.54	0.42
5:H:182:LEU:HB2	5:H:185:SER:HB3	2.01	0.42
5:H:10:TYR:HB3	5:H:153:HIS:ND1	2.34	0.42
4:I:9:PRO:O	4:I:9:PRO:CD	2.66	0.42
1:A:7:TYR:O	1:A:98:MET:HA	2.19	0.42
2:B:55:SER:OG	2:B:56:PHE:N	2.52	0.42
2:E:41:LYS:HG3	2:E:78:TYR:CE1	2.55	0.42
1:D:194:VAL:HG23	1:D:195:SER:N	2.35	0.42
4:I:96:ASP:OD2	5:J:100:VAL:HG21	2.20	0.42
5:H:30:GLU:HG3	5:H:96:LEU:CB	2.49	0.42
5:H:10:TYR:HB3	5:H:153:HIS:HD1	1.85	0.42
1:A:156:LEU:HD13	3:C:5:GLY:HA3	2.02	0.42
5:J:14:ARG:O	5:J:17:GLU:HB2	2.20	0.42
1:A:197:HIS:C	1:A:251:SER:HB2	2.41	0.41
1:A:14:ARG:NH2	1:A:39:ASP:OD1	2.43	0.41
1:D:137:ASP:OD1	1:D:137:ASP:C	2.58	0.41
1:D:98:MET:HG3	1:D:98:MET:O	2.19	0.41
5:J:169:VAL:HA	5:J:192:ARG:O	2.20	0.41
2:E:59:ASP:C	2:E:59:ASP:OD1	2.58	0.41
1:D:254:GLU:HG2	1:D:274:TRP:NE1	2.34	0.41
1:A:209:TYR:CG	1:A:210:PRO:HA	2.55	0.41
4:I:124:ASP:OD1	4:I:125:SER:N	2.53	0.41
4:I:42:SER:HB2	5:J:90:LEU:CD2	2.51	0.41
2:E:23:LEU:O	2:E:67:TYR:HA	2.20	0.41
5:J:113:VAL:HG22	5:J:118:LYS:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:17:ASN:OD1	2:B:97:ARG:NH2	2.51	0.41
4:G:35:PHE:HB2	4:G:91:ALA:HB3	2.01	0.41
1:A:170:ARG:HH12	4:I:27:ARG:HH22	1.69	0.41
5:H:133:GLU:HG2	5:H:139:LYS:O	2.20	0.41
2:B:17:ASN:ND2	2:B:72:PRO:CB	2.64	0.41
5:J:101:GLU:H	5:J:101:GLU:CD	2.24	0.41
1:A:204:TRP:CH2	2:B:99:MET:C	2.93	0.41
1:A:217:TRP:O	1:A:218:GLN:NE2	2.54	0.41
5:J:182:LEU:HD23	5:J:184:ASP:C	2.41	0.41
4:I:117:PRO:HB2	4:I:196:THR:HB	2.02	0.41
5:H:25:GLN:O	5:H:72:GLU:HB2	2.20	0.41
1:A:123:TYR:HH	1:A:143:THR:HG1	1.62	0.41
1:D:266:LEU:HD22	1:D:270:LEU:HD13	2.02	0.40
3:F:7:LEU:HD11	5:H:95:PHE:HB2	2.03	0.40
5:H:119:VAL:HG13	5:H:150:TYR:O	2.21	0.40
1:A:195:SER:OG	1:A:196:ASP:N	2.48	0.40
1:D:11:SER:HA	1:D:21:ARG:O	2.21	0.40
5:H:44:ARG:HD2	5:H:60:ILE:HD12	2.03	0.40
1:D:123:TYR:O	1:D:140:ALA:HB1	2.21	0.40
1:D:218:GLN:HA	1:D:218:GLN:HE21	1.87	0.40
4:I:173:VAL:HG12	5:J:192:ARG:HH22	1.86	0.40
1:D:106:ASP:OD1	1:D:108:ARG:NH1	2.55	0.40
5:J:148:GLY:HA2	5:J:186:ARG:HB3	2.03	0.40
2:E:39:LEU:HD13	2:E:68:THR:HG22	2.03	0.40
1:A:258:THR:HG21	1:A:260:HIS:CE1	2.57	0.40
4:I:124:ASP:C	4:I:124:ASP:OD1	2.60	0.40
5:J:172:ASP:OD2	5:J:190:SER:OG	2.38	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	263/275 (96%)	242 (92%)	19 (7%)	2 (1%)	24	58
1	D	261/275 (95%)	245 (94%)	16 (6%)	0	100	100
2	B	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	E	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
3	C	7/9 (78%)	5 (71%)	2 (29%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	G	199/206 (97%)	179 (90%)	19 (10%)	1 (0%)	34	69
4	I	199/206 (97%)	187 (94%)	12 (6%)	0	100	100
5	H	239/243 (98%)	228 (95%)	10 (4%)	1 (0%)	39	73
5	J	238/243 (98%)	221 (93%)	17 (7%)	0	100	100
All	All	1609/1666 (97%)	1501 (93%)	104 (6%)	4 (0%)	52	83

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	SER
5	H	219	ALA
1	A	251	SER
4	G	95	GLY

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	224/231 (97%)	221 (99%)	3 (1%)	76	93
1	D	222/231 (96%)	220 (99%)	2 (1%)	84	95
2	B	95/95 (100%)	94 (99%)	1 (1%)	80	94
2	E	95/95 (100%)	95 (100%)	0	100	100
3	C	6/6 (100%)	6 (100%)	0	100	100
3	F	6/6 (100%)	6 (100%)	0	100	100
4	G	174/179 (97%)	171 (98%)	3 (2%)	68	90

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	I	174/179 (97%)	173 (99%)	1 (1%)	90	97
5	H	212/214 (99%)	212 (100%)	0	100	100
5	J	211/214 (99%)	210 (100%)	1 (0%)	92	97
All	All	1419/1450 (98%)	1408 (99%)	11 (1%)	86	96

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	113	TYR
1	D	218	GLN
1	A	156	LEU
1	A	177	GLU
1	A	197	HIS
2	B	85	VAL
4	G	70	SER
4	G	73	VAL
4	G	143	GLN
4	I	132	VAL
5	J	72	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	3	HIS
1	D	70	HIS
1	D	114	HIS
1	D	197	HIS
1	D	218	GLN
1	D	255	GLN
2	E	2	GLN
1	A	3	HIS
1	A	70	HIS
1	A	114	HIS
1	A	260	HIS
2	B	2	GLN
2	B	17	ASN
4	G	7	GLN
4	G	172	ASN
4	I	21	ASN
4	I	52	ASN

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Mol	Chain	Res	Type
4	I	69	GLN
4	I	111	ASN
4	I	170	ASN
4	I	188	ASN
5	H	25	GLN
5	H	174	GLN
5	H	201	GLN
5	J	25	GLN
5	J	201	GLN
5	J	212	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 carbohydrates 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	267/275 (97%)	-0.14	7 (2%) 59 55	20, 37, 67, 84	5 (1%)
1	D	265/275 (96%)	-0.09	12 (4%) 37 31	19, 35, 74, 91	5 (1%)
2	B	100/100 (100%)	-0.20	2 (2%) 68 65	21, 34, 54, 75	1 (1%)
2	E	100/100 (100%)	-0.24	1 (1%) 84 82	21, 33, 53, 72	1 (1%)
3	C	9/9 (100%)	-0.48	0 100 100	36, 38, 38, 39	0
3	F	9/9 (100%)	-0.50	0 100 100	29, 31, 33, 34	0
4	G	201/206 (97%)	0.65	22 (10%) 7 4	39, 63, 98, 107	2 (0%)
4	I	201/206 (97%)	0.85	31 (15%) 3 1	34, 65, 119, 144	3 (1%)
5	H	241/243 (99%)	0.32	10 (4%) 41 35	39, 56, 77, 91	0
5	J	240/243 (98%)	0.44	10 (4%) 40 34	40, 61, 93, 115	0
All	All	1633/1666 (98%)	0.22	95 (5%) 26 21	19, 50, 93, 144	17 (1%)

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	I	201	PRO	9.8
4	I	127	SER	9.3
5	J	225	ASP	5.3
4	G	203	PRO	5.0
4	I	200	SER	4.8
4	I	129	ASP	4.8
1	D	274	TRP	4.6
5	J	135	SER	4.4
4	I	180	ASP	4.2
4	I	165	MET	4.0
2	B	0	MET	4.0
4	I	96	ASP	4.0
5	H	223	THR	4.0

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Mol	Chain	Res	Type	RSRZ
4	I	195	ASP	3.7
4	G	181	SER	3.7
1	D	273	ARG	3.5
4	G	165	ARG	3.5
4	I	94	GLY	3.5
5	J	221	GLU	3.4
4	I	179	SER	3.4
4	I	1	LYS	3.4
4	G	166	SER	3.4
4	I	164	SER	3.4
4	I	95	ALA	3.3
4	I	184	ALA	3.3
1	A	274	TRP	3.2
4	G	129	SER	3.2
2	E	0	MET	3.2
4	G	182	ASP	3.2
5	H	26	ASP	3.2
4	G	150	LYS	3.1
4	I	133	CYS	3.1
4	I	148	LYS	3.1
5	H	40	GLY	3.0
1	D	17	ARG	3.0
5	H	135	SER	3.0
4	I	102	LYS	3.0
5	H	243	ASP	3.0
5	J	241	ARG	2.9
4	I	163	ARG	2.9
4	G	132	LYS	2.9
4	G	96	GLY	2.9
4	I	128	ALA	2.9
5	H	180	PRO	2.9
5	J	183	ASN	2.9
1	A	137	ASP	2.8
4	I	178	LYS	2.8
5	H	225	ASP	2.8
4	I	136	THR	2.8
4	I	196	THR	2.8
5	J	180	PRO	2.7
4	G	9	SER	2.7
4	G	167	MET	2.7
1	D	218	GLN	2.7
4	I	131	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	196	ASP	2.6
1	D	256	ARG	2.6
4	I	150	SER	2.6
4	G	3	LYS	2.6
1	A	219	ARG	2.5
4	G	152	SER	2.5
4	I	92	GLY	2.5
4	G	98	ASP	2.5
1	D	138	MET	2.5
4	G	138	THR	2.5
4	G	82	PRO	2.5
5	H	92	ALA	2.5
4	I	126	LYS	2.4
4	G	196	GLU	2.4
4	I	188	ASN	2.4
4	I	189	ASN	2.4
1	D	194	VAL	2.4
1	D	254	GLU	2.3
4	G	191	ASN	2.3
5	H	138	GLN	2.3
2	B	48	LYS	2.3
4	G	197	ASP	2.2
1	A	203	CYS	2.2
1	D	250	PRO	2.2
5	H	91	CYS	2.2
5	J	52	VAL	2.2
1	D	253	GLN	2.2
4	I	134	LEU	2.2
4	G	148	GLN	2.2
4	I	194	GLU	2.1
1	A	218	GLN	2.1
4	G	183	PHE	2.1
1	A	267	PRO	2.1
4	G	190	ASN	2.0
1	D	195	SER	2.0
5	J	53	LYS	2.0
4	I	186	ALA	2.0
5	J	20	PHE	2.0
5	J	82	SER	2.0
1	D	228	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.