



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:06 AM UTC

PDB ID : 4FRW / pdb_00004frw
Title : Crystal structure of human nectin-4 extracellular fragment D1-D2
Authors : Harrison, O.J.; Jin, X.; Brasch, J.; Shapiro, L.
Deposited on : 2012-06-26
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 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

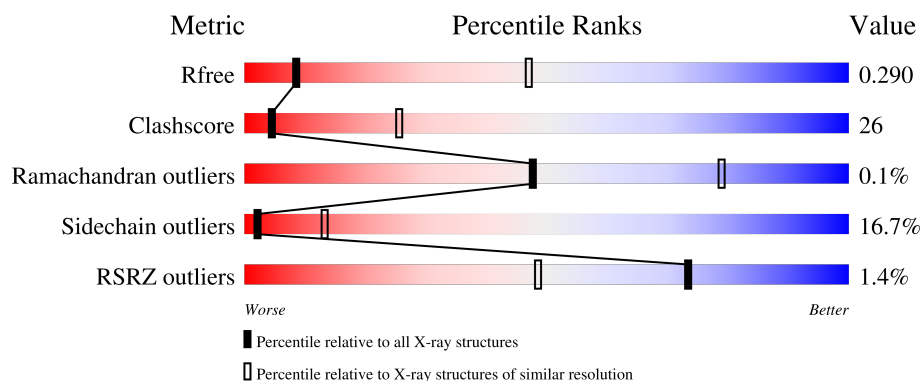
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	218	
1	B	218	
1	C	218	
1	D	218	
1	E	218	

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Mol	Chain	Length	Quality of chain
1	F	218	6% 41% 36% 10% 13%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9257 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 Poliovirus receptor-related protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1611	1002	291	313	5			
1	B	213	Total	C	N	O	S	0	0	0
			1601	996	289	311	5			
1	C	211	Total	C	N	O	S	0	0	0
			1571	978	280	308	5			
1	D	212	Total	C	N	O	S	0	0	0
			1561	973	278	305	5			
1	E	208	Total	C	N	O	S	0	0	0
			1533	954	276	298	5			
1	F	189	Total	C	N	O	S	0	0	0
			1380	860	246	269	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	HIS	-	expression tag	UNP Q96NY8
A	245	HIS	-	expression tag	UNP Q96NY8
A	246	HIS	-	expression tag	UNP Q96NY8
A	247	HIS	-	expression tag	UNP Q96NY8
A	248	HIS	-	expression tag	UNP Q96NY8
A	249	HIS	-	expression tag	UNP Q96NY8
B	244	HIS	-	expression tag	UNP Q96NY8
B	245	HIS	-	expression tag	UNP Q96NY8
B	246	HIS	-	expression tag	UNP Q96NY8
B	247	HIS	-	expression tag	UNP Q96NY8
B	248	HIS	-	expression tag	UNP Q96NY8
B	249	HIS	-	expression tag	UNP Q96NY8
C	244	HIS	-	expression tag	UNP Q96NY8
C	245	HIS	-	expression tag	UNP Q96NY8
C	246	HIS	-	expression tag	UNP Q96NY8
C	247	HIS	-	expression tag	UNP Q96NY8
C	248	HIS	-	expression tag	UNP Q96NY8

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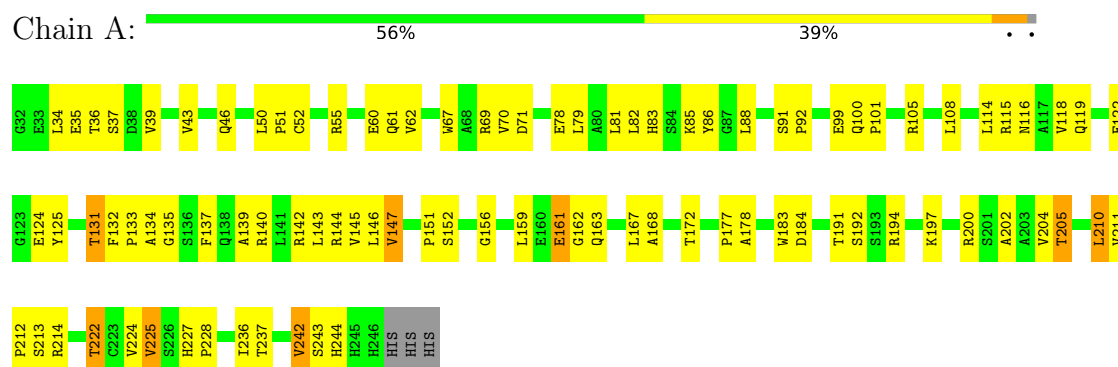
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Chain	Residue	Modelled	Actual	Comment	Reference
C	249	HIS	-	expression tag	UNP Q96NY8
D	244	HIS	-	expression tag	UNP Q96NY8
D	245	HIS	-	expression tag	UNP Q96NY8
D	246	HIS	-	expression tag	UNP Q96NY8
D	247	HIS	-	expression tag	UNP Q96NY8
D	248	HIS	-	expression tag	UNP Q96NY8
D	249	HIS	-	expression tag	UNP Q96NY8
E	244	HIS	-	expression tag	UNP Q96NY8
E	245	HIS	-	expression tag	UNP Q96NY8
E	246	HIS	-	expression tag	UNP Q96NY8
E	247	HIS	-	expression tag	UNP Q96NY8
E	248	HIS	-	expression tag	UNP Q96NY8
E	249	HIS	-	expression tag	UNP Q96NY8
F	244	HIS	-	expression tag	UNP Q96NY8
F	245	HIS	-	expression tag	UNP Q96NY8
F	246	HIS	-	expression tag	UNP Q96NY8
F	247	HIS	-	expression tag	UNP Q96NY8
F	248	HIS	-	expression tag	UNP Q96NY8
F	249	HIS	-	expression tag	UNP Q96NY8

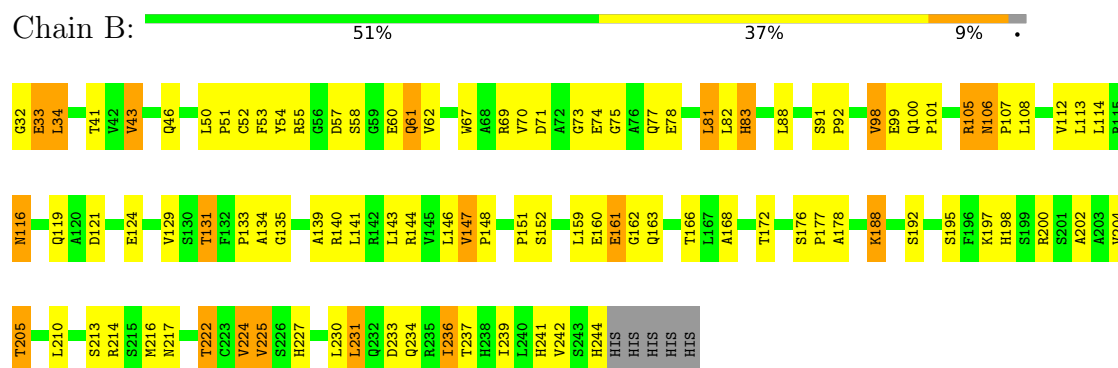
3 Residue-property plots [i](#)

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.

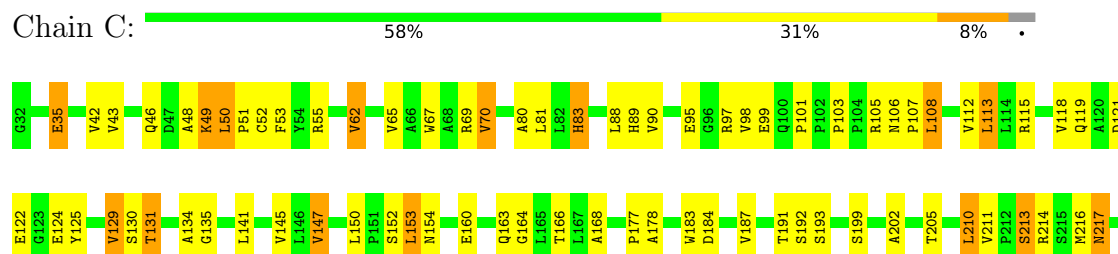
• Molecule 1: Poliovirus receptor-related protein 4



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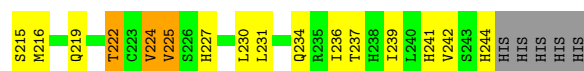


• Molecule 1: Poliovirus receptor-related protein 4

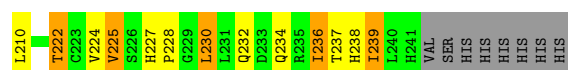




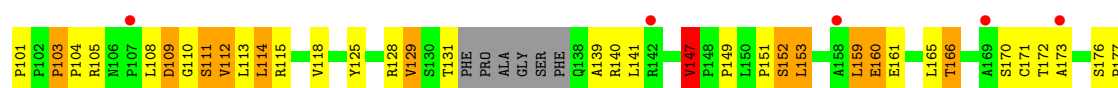
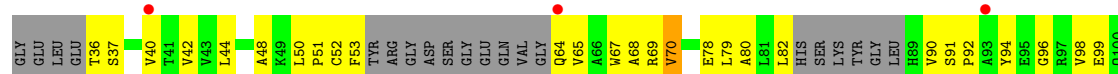
• Molecule 1: Poliovirus receptor-related protein 4



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• Molecule 1: Poliovirus receptor-related protein 4



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	85.97Å 142.81Å 341.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.50 20.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (20.00-3.50) 98.7 (20.00-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 3.52Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.257 , 0.290 0.260 , 0.290	Depositor DCC
R_{free} test set	1334 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	96.9	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 58.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.054 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.055 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	9257	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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	1.00	1/1649 (0.1%)	1.09	5/2251 (0.2%)
1	B	1.00	2/1639 (0.1%)	1.03	5/2237 (0.2%)
1	C	1.00	3/1608 (0.2%)	1.07	8/2199 (0.4%)
1	D	0.96	1/1598 (0.1%)	1.03	1/2187 (0.0%)
1	E	0.98	0/1569	1.12	5/2148 (0.2%)
1	F	0.99	0/1409	1.14	6/1930 (0.3%)
All	All	0.99	7/9472 (0.1%)	1.08	30/12952 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	83	HIS	CG-ND1	-5.95	1.31	1.38
1	D	241	HIS	CG-CD2	5.75	1.42	1.35
1	C	83	HIS	CA-C	-5.64	1.46	1.52
1	B	83	HIS	CG-ND1	-5.60	1.32	1.38
1	C	89	HIS	CG-CD2	5.19	1.41	1.35
1	C	241	HIS	CG-CD2	5.13	1.41	1.35
1	B	198	HIS	CG-ND1	-5.07	1.32	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	50	LEU	CA-C-N	9.14	129.50	119.90
1	A	50	LEU	C-N-CA	9.14	129.50	119.90
1	E	87	GLY	N-CA-C	8.41	123.45	110.91
1	C	147	VAL	CA-C-N	6.76	124.53	119.66
1	C	147	VAL	C-N-CA	6.76	124.53	119.66
1	B	74	GLU	N-CA-C	-6.58	105.13	113.02
1	A	147	VAL	CA-C-N	6.39	124.26	119.66
1	A	147	VAL	C-N-CA	6.39	124.26	119.66
1	F	70	VAL	N-CA-C	6.27	118.51	108.85
1	A	162	GLY	N-CA-C	6.25	120.05	112.49
1	E	163	GLN	N-CA-C	6.03	117.30	107.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	147	VAL	CA-C-N	5.98	123.97	119.66
1	E	147	VAL	C-N-CA	5.98	123.97	119.66
1	D	163	GLN	N-CA-C	-5.90	106.32	112.93
1	F	147	VAL	CA-C-N	5.68	123.81	119.66
1	F	147	VAL	C-N-CA	5.68	123.81	119.66
1	F	244	HIS	N-CA-C	5.52	117.58	108.48
1	C	214	ARG	CA-C-N	5.48	130.37	121.98
1	C	214	ARG	C-N-CA	5.48	130.37	121.98
1	C	236	ILE	CB-CA-C	-5.48	102.66	110.83
1	B	227	HIS	CA-C-N	-5.47	114.27	120.89
1	B	227	HIS	C-N-CA	-5.47	114.27	120.89
1	F	103	PRO	N-CA-C	5.46	117.36	110.70
1	E	237	THR	N-CA-C	5.23	116.93	108.41
1	C	90	VAL	N-CA-CB	5.19	116.38	110.72
1	F	128	ARG	N-CA-C	5.17	116.84	108.26
1	B	106	ASN	CA-C-N	5.12	124.83	119.82
1	B	106	ASN	C-N-CA	5.12	124.83	119.82
1	C	227	HIS	CA-C-N	5.11	124.61	119.19
1	C	227	HIS	C-N-CA	5.11	124.61	119.19

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	1611	0	1566	77	0
1	B	1601	0	1562	65	0
1	C	1571	0	1517	65	0
1	D	1561	0	1505	66	0
1	E	1533	0	1478	87	0
1	F	1380	0	1328	122	0
All	All	9257	0	8956	478	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 26.

All (478) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:66:ALA:HB1	1:E:128:ARG:CZ	1.22	1.62
1:E:66:ALA:CB	1:E:128:ARG:CZ	2.00	1.40
1:E:66:ALA:HB1	1:E:128:ARG:NH2	1.43	1.33
1:E:101:PRO:O	1:E:105:ARG:HD3	1.38	1.19
1:E:66:ALA:CB	1:E:128:ARG:NE	2.06	1.18
1:F:153:LEU:HD23	1:F:236:ILE:HG22	1.26	1.17
1:A:131:THR:CG2	1:A:135:GLY:O	2.01	1.08
1:E:79:LEU:HG	1:E:94:TYR:CE1	1.89	1.07
1:F:51:PRO:HB3	1:F:111:SER:OG	1.55	1.06
1:F:151:PRO:HA	1:F:172:THR:O	1.56	1.05
1:F:50:LEU:HD22	1:F:141:LEU:CD1	1.88	1.03
1:A:70:VAL:HG11	1:A:140:ARG:NH1	1.72	1.03
1:E:66:ALA:HB3	1:E:128:ARG:CG	1.88	1.02
1:F:50:LEU:HD22	1:F:141:LEU:HD13	1.38	1.02
1:E:101:PRO:O	1:E:105:ARG:NH1	1.94	1.01
1:A:131:THR:HG23	1:A:135:GLY:O	1.62	0.98
1:C:101:PRO:O	1:C:105:ARG:HD3	1.63	0.97
1:F:40:VAL:HG21	1:F:50:LEU:HD23	1.45	0.97
1:F:153:LEU:HD23	1:F:236:ILE:CG2	1.95	0.96
1:E:66:ALA:HB3	1:E:128:ARG:NE	1.76	0.95
1:E:41:THR:HG23	1:E:144:ARG:HG3	1.50	0.92
1:E:81:LEU:O	1:E:88:LEU:HG	1.67	0.92
1:A:131:THR:HG22	1:A:135:GLY:O	1.68	0.92
1:F:153:LEU:CD2	1:F:236:ILE:HG22	1.99	0.92
1:F:191:THR:HB	1:F:207:GLU:O	1.70	0.92
1:F:53:PHE:HA	1:F:109:ASP:OD2	1.71	0.91
1:A:101:PRO:O	1:A:105:ARG:HD3	1.69	0.91
1:F:183:TRP:CD1	1:F:191:THR:HG21	2.09	0.88
1:E:101:PRO:O	1:E:105:ARG:CD	2.21	0.88
1:C:213:SER:O	1:C:242:VAL:HG21	1.74	0.88
1:F:51:PRO:CB	1:F:111:SER:OG	2.22	0.87
1:A:70:VAL:HG11	1:A:140:ARG:HH11	1.35	0.87
1:F:53:PHE:HA	1:F:109:ASP:CG	2.00	0.87
1:B:168:ALA:HB2	1:B:210:LEU:HD22	1.58	0.86
1:E:79:LEU:HG	1:E:94:TYR:CZ	2.11	0.85
1:D:120:ALA:HB2	1:D:200:ARG:HH12	1.40	0.85
1:A:70:VAL:HG21	1:A:124:GLU:CB	2.08	0.84
1:C:160:GLU:HB3	1:C:163:GLN:HE22	1.42	0.84
1:A:70:VAL:HG21	1:A:124:GLU:HB3	1.62	0.81
1:B:78:GLU:OE2	1:B:81:LEU:HD23	1.81	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:54:TYR:CZ	1:E:131:THR:HG21	2.16	0.80
1:E:60:GLU:HB2	1:E:132:PHE:O	1.80	0.80
1:F:176:SER:HB2	1:F:201:SER:HA	1.63	0.80
1:B:217:ASN:ND2	1:B:241:HIS:HA	1.95	0.80
1:C:49:LYS:HA	1:C:113:LEU:HD12	1.63	0.79
1:A:91:SER:OG	1:A:92:PRO:HD2	1.81	0.79
1:A:194:ARG:NH1	1:B:148:PRO:O	2.14	0.79
1:A:70:VAL:CG1	1:A:140:ARG:NH1	2.45	0.79
1:F:151:PRO:HG3	1:F:173:ALA:HB2	1.65	0.79
1:A:172:THR:HG23	1:A:205:THR:HG22	1.66	0.78
1:E:54:TYR:CE1	1:E:131:THR:HG21	2.18	0.78
1:F:185:THR:HG22	1:F:186:GLU:H	1.48	0.78
1:B:70:VAL:CG2	1:B:124:GLU:HB3	2.13	0.78
1:F:151:PRO:HG2	1:F:225:VAL:HG11	1.66	0.77
1:E:66:ALA:CB	1:E:128:ARG:NH2	2.31	0.77
1:D:216:MET:HA	1:D:219:GLN:HG3	1.65	0.77
1:F:53:PHE:CE2	1:F:108:LEU:HB3	2.20	0.76
1:F:105:ARG:HH12	1:F:110:GLY:HA3	1.51	0.76
1:E:66:ALA:HB3	1:E:128:ARG:HG3	1.64	0.75
1:F:51:PRO:CA	1:F:111:SER:OG	2.34	0.75
1:F:82:LEU:HD21	1:F:112:VAL:HG23	1.67	0.74
1:F:194:ARG:CG	1:F:194:ARG:HH11	1.99	0.74
1:B:222:THR:HG22	1:B:237:THR:HG23	1.68	0.74
1:E:90:VAL:HG21	1:E:94:TYR:O	1.87	0.74
1:B:147:VAL:HG13	1:B:176:SER:HB3	1.68	0.74
1:F:53:PHE:HD2	1:F:108:LEU:HD23	1.51	0.74
1:B:43:VAL:HG22	1:B:46:GLN:HG2	1.69	0.73
1:B:230:LEU:HD21	1:B:233:ASP:HA	1.70	0.73
1:A:70:VAL:HG22	1:A:124:GLU:C	2.13	0.73
1:D:121:ASP:OD1	1:D:145:VAL:HG21	1.87	0.73
1:F:80:ALA:HB3	1:F:98:VAL:HG11	1.71	0.73
1:E:60:GLU:CB	1:E:132:PHE:O	2.37	0.73
1:F:165:LEU:HB3	1:F:211:VAL:HG13	1.70	0.73
1:F:50:LEU:HD22	1:F:141:LEU:HD11	1.71	0.72
1:E:66:ALA:HB3	1:E:128:ARG:CD	2.19	0.72
1:A:70:VAL:CG2	1:A:124:GLU:CB	2.67	0.72
1:F:101:PRO:O	1:F:105:ARG:HD3	1.88	0.72
1:D:161:GLU:OE2	1:D:161:GLU:HA	1.88	0.72
1:E:101:PRO:C	1:E:105:ARG:HD3	2.13	0.72
1:E:64:GLN:HE21	1:E:81:LEU:HD21	1.53	0.72
1:B:69:ARG:HG2	1:B:77:GLN:HB3	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:232:GLN:HA	1:E:232:GLN:NE2	2.05	0.71
1:C:70:VAL:CG2	1:C:124:GLU:HB3	2.20	0.71
1:D:65:VAL:HG22	1:D:129:VAL:HG23	1.73	0.71
1:C:131:THR:HG23	1:C:135:GLY:O	1.91	0.71
1:E:224:VAL:HG23	1:E:224:VAL:O	1.91	0.70
1:F:153:LEU:CD2	1:F:236:ILE:CG2	2.62	0.70
1:D:159:LEU:O	1:D:242:VAL:HA	1.91	0.70
1:E:41:THR:HG23	1:E:144:ARG:CG	2.22	0.70
1:F:194:ARG:HH11	1:F:194:ARG:HG2	1.56	0.70
1:E:65:VAL:O	1:E:81:LEU:HD12	1.91	0.69
1:D:54:TYR:CE2	1:D:131:THR:HG21	2.28	0.69
1:F:183:TRP:HD1	1:F:191:THR:HG21	1.55	0.69
1:E:51:PRO:HD2	1:E:141:LEU:HD22	1.73	0.69
1:F:109:ASP:OD2	1:F:109:ASP:C	2.36	0.69
1:C:65:VAL:HG22	1:C:129:VAL:HG22	1.76	0.68
1:D:69:ARG:NH1	1:D:94:TYR:OH	2.27	0.68
1:A:70:VAL:CG1	1:A:140:ARG:HH12	2.05	0.68
1:F:151:PRO:CA	1:F:172:THR:O	2.39	0.68
1:F:103:PRO:HA	1:F:105:ARG:HG3	1.76	0.68
1:F:53:PHE:CD2	1:F:108:LEU:HD23	2.29	0.67
1:C:70:VAL:HG23	1:C:124:GLU:HB3	1.74	0.67
1:E:66:ALA:HB3	1:E:128:ARG:HG2	1.76	0.66
1:E:178:ALA:HB2	1:E:202:ALA:HB3	1.77	0.66
1:C:98:VAL:HG13	1:C:112:VAL:CG1	2.26	0.66
1:A:70:VAL:CG2	1:A:124:GLU:HB2	2.26	0.66
1:A:168:ALA:HB2	1:A:210:LEU:HD22	1.76	0.66
1:C:50:LEU:N	1:C:50:LEU:CD1	2.60	0.65
1:A:70:VAL:CG2	1:A:124:GLU:C	2.69	0.65
1:A:70:VAL:HG21	1:A:124:GLU:HB2	1.79	0.65
1:B:131:THR:OG1	1:B:134:ALA:HB3	1.96	0.65
1:F:191:THR:CB	1:F:207:GLU:O	2.44	0.65
1:B:62:VAL:HG13	1:B:108:LEU:HG	1.77	0.65
1:F:192:SER:OG	1:F:207:GLU:HB2	1.97	0.64
1:B:34:LEU:HD13	1:B:139:ALA:HB2	1.79	0.64
1:E:81:LEU:O	1:E:88:LEU:CG	2.42	0.64
1:C:80:ALA:HB1	1:C:88:LEU:HD11	1.79	0.64
1:F:50:LEU:CD2	1:F:141:LEU:HD13	2.20	0.64
1:F:37:SER:HB2	1:F:51:PRO:HD2	1.80	0.64
1:A:39:VAL:HG23	1:A:142:ARG:O	1.98	0.64
1:D:69:ARG:NH2	1:D:120:ALA:O	2.31	0.63
1:F:94:TYR:HB3	1:F:98:VAL:HG23	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:53:PHE:CA	1:F:109:ASP:CG	2.72	0.63
1:F:80:ALA:HB2	1:F:90:VAL:HG23	1.80	0.63
1:C:50:LEU:N	1:C:50:LEU:HD12	2.13	0.63
1:F:187:VAL:HG23	1:F:208:PHE:CZ	2.34	0.63
1:D:174:GLU:HG3	1:D:202:ALA:O	1.99	0.63
1:F:151:PRO:CG	1:F:173:ALA:HB2	2.29	0.62
1:D:116:ASN:O	1:D:118:VAL:HG23	1.99	0.62
1:D:118:VAL:HG12	1:D:119:GLN:N	2.14	0.62
1:F:40:VAL:HG21	1:F:50:LEU:CD2	2.25	0.62
1:F:178:ALA:HB2	1:F:202:ALA:HB3	1.82	0.62
1:E:153:LEU:HD21	1:E:238:HIS:HB2	1.82	0.62
1:E:91:SER:HB3	1:E:92:PRO:HD2	1.81	0.62
1:C:217:ASN:HD21	1:C:242:VAL:HB	1.63	0.62
1:A:52:CYS:HB2	1:A:67:TRP:CZ2	2.35	0.61
1:B:168:ALA:HB2	1:B:210:LEU:CD2	2.29	0.61
1:B:70:VAL:HG23	1:B:124:GLU:HB3	1.81	0.61
1:B:172:THR:HG23	1:B:205:THR:HG22	1.82	0.61
1:A:82:LEU:HB3	1:A:88:LEU:HD12	1.83	0.61
1:F:80:ALA:CB	1:F:98:VAL:HG11	2.31	0.61
1:F:42:VAL:HG21	1:F:48:ALA:HB2	1.82	0.61
1:F:51:PRO:HA	1:F:111:SER:OG	2.01	0.61
1:B:101:PRO:O	1:B:105:ARG:NH1	2.33	0.61
1:A:99:GLU:OE2	1:A:115:ARG:HD2	2.01	0.60
1:F:79:LEU:HD22	1:F:94:TYR:CZ	2.36	0.60
1:C:99:GLU:OE1	1:C:115:ARG:HD2	2.01	0.60
1:A:70:VAL:HG22	1:A:124:GLU:O	2.02	0.60
1:D:184:ASP:O	1:D:222:THR:HG23	2.01	0.60
1:A:34:LEU:HD23	1:A:139:ALA:HB2	1.83	0.60
1:E:232:GLN:HA	1:E:232:GLN:HE21	1.64	0.60
1:F:191:THR:HA	1:F:207:GLU:O	2.01	0.60
1:A:43:VAL:HG23	1:A:46:GLN:HG2	1.83	0.60
1:D:43:VAL:HG22	1:D:46:GLN:HG2	1.82	0.60
1:E:61:GLN:HE22	1:E:107:PRO:HB2	1.65	0.60
1:A:79:LEU:O	1:A:91:SER:HB3	2.02	0.60
1:F:151:PRO:HG3	1:F:173:ALA:CB	2.32	0.60
1:E:98:VAL:HG13	1:E:112:VAL:CG1	2.32	0.59
1:E:98:VAL:HG13	1:E:112:VAL:HG11	1.83	0.59
1:E:34:LEU:HD21	1:E:129:VAL:HG11	1.85	0.59
1:C:51:PRO:HD2	1:C:141:LEU:HD22	1.84	0.59
1:B:53:PHE:CE2	1:B:55:ARG:HD2	2.37	0.59
1:C:106:ASN:OD1	1:C:108:LEU:CD1	2.51	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:VAL:HG12	1:C:119:GLN:N	2.18	0.59
1:F:50:LEU:CD2	1:F:141:LEU:CD1	2.74	0.59
1:F:210:LEU:HD13	1:F:216:MET:HE1	1.84	0.59
1:C:98:VAL:CG1	1:C:112:VAL:CG1	2.81	0.59
1:F:64:GLN:O	1:F:129:VAL:CG2	2.51	0.58
1:E:230:LEU:HD21	1:E:234:GLN:NE2	2.18	0.58
1:F:80:ALA:HB2	1:F:98:VAL:HG21	1.84	0.58
1:D:54:TYR:O	1:D:108:LEU:HD23	2.04	0.58
1:B:61:GLN:NE2	1:B:62:VAL:H	2.02	0.58
1:B:178:ALA:HB2	1:B:202:ALA:HB3	1.85	0.58
1:D:119:GLN:O	1:D:120:ALA:HB3	2.03	0.58
1:D:161:GLU:OE1	1:D:214:ARG:HG2	2.04	0.58
1:B:121:ASP:O	1:B:143:LEU:HD23	2.04	0.58
1:F:183:TRP:CD1	1:F:191:THR:CG2	2.85	0.58
1:B:50:LEU:N	1:B:50:LEU:HD12	2.18	0.58
1:F:40:VAL:CG2	1:F:50:LEU:HD23	2.29	0.57
1:F:50:LEU:HD13	1:F:141:LEU:HD11	1.86	0.57
1:A:227:HIS:CG	1:A:228:PRO:HD2	2.38	0.57
1:B:98:VAL:HG22	1:B:112:VAL:CG1	2.34	0.57
1:E:118:VAL:HG12	1:E:119:GLN:N	2.19	0.57
1:B:160:GLU:HG2	1:B:162:GLY:H	1.69	0.57
1:B:231:LEU:CD2	1:B:231:LEU:H	2.18	0.57
1:D:69:ARG:HB3	1:D:77:GLN:HB3	1.85	0.57
1:E:61:GLN:NE2	1:E:107:PRO:HB2	2.20	0.57
1:D:187:VAL:HB	1:D:216:MET:HE2	1.87	0.57
1:E:43:VAL:HG13	1:E:46:GLN:HG3	1.86	0.57
1:B:50:LEU:N	1:B:50:LEU:CD1	2.68	0.57
1:E:64:GLN:HG2	1:E:81:LEU:HD11	1.87	0.57
1:F:227:HIS:CE1	1:F:228:PRO:O	2.58	0.57
1:B:131:THR:HG23	1:B:135:GLY:O	2.05	0.56
1:D:100:GLN:CG	1:D:105:ARG:HD3	2.35	0.56
1:F:103:PRO:CB	1:F:104:PRO:HA	2.35	0.56
1:F:194:ARG:O	1:F:204:VAL:HG23	2.04	0.56
1:F:231:LEU:H	1:F:231:LEU:HD12	1.69	0.56
1:E:151:PRO:HG3	1:E:225:VAL:HG11	1.86	0.56
1:F:79:LEU:HD22	1:F:94:TYR:CE2	2.40	0.56
1:F:194:ARG:HD3	1:F:195:SER:H	1.69	0.56
1:D:34:LEU:HD21	1:D:52:CYS:SG	2.46	0.56
1:A:184:ASP:O	1:A:222:THR:HG23	2.05	0.56
1:B:163:GLN:HG3	1:B:166:THR:HG21	1.87	0.56
1:C:106:ASN:OD1	1:C:107:PRO:HD2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:185:THR:HG22	1:F:186:GLU:N	2.19	0.56
1:A:37:SER:HB2	1:A:51:PRO:HG2	1.88	0.56
1:C:141:LEU:C	1:C:141:LEU:HD12	2.31	0.56
1:F:149:PRO:HB2	1:F:173:ALA:HB1	1.87	0.56
1:D:210:LEU:O	1:D:210:LEU:HD23	2.06	0.55
1:B:146:LEU:HD22	1:B:177:PRO:HD3	1.88	0.55
1:D:214:ARG:NH2	1:D:244:HIS:CD2	2.74	0.55
1:B:161:GLU:OE1	1:B:214:ARG:NH1	2.39	0.55
1:B:82:LEU:HD12	1:B:82:LEU:O	2.07	0.55
1:D:69:ARG:HH12	1:D:94:TYR:HH	1.54	0.55
1:D:131:THR:C	1:D:132:PHE:CD2	2.85	0.54
1:F:160:GLU:O	1:F:212:PRO:HG2	2.07	0.54
1:C:35:GLU:OE1	1:C:55:ARG:NH1	2.40	0.54
1:E:159:LEU:HD22	1:E:166:THR:OG1	2.07	0.54
1:A:43:VAL:CG2	1:A:46:GLN:HG2	2.38	0.54
1:B:112:VAL:HG12	1:B:113:LEU:N	2.23	0.54
1:C:217:ASN:ND2	1:C:242:VAL:HB	2.23	0.54
1:A:214:ARG:NH1	1:A:242:VAL:O	2.39	0.54
1:D:91:SER:HB2	1:D:92:PRO:HD2	1.88	0.54
1:B:151:PRO:HB3	1:B:225:VAL:HG11	1.90	0.54
1:D:161:GLU:OE2	1:D:161:GLU:CA	2.54	0.54
1:E:66:ALA:HB1	1:E:128:ARG:NH1	2.07	0.54
1:D:106:ASN:OD1	1:D:107:PRO:HD2	2.08	0.54
1:D:230:LEU:HD21	1:D:234:GLN:HG2	1.91	0.54
1:C:122:GLU:HG3	1:C:145:VAL:H	1.73	0.53
1:C:222:THR:HB	1:C:237:THR:OG1	2.08	0.53
1:E:153:LEU:HD12	1:E:170:SER:O	2.08	0.53
1:F:82:LEU:HD12	1:F:82:LEU:O	2.07	0.53
1:A:131:THR:CG2	1:A:134:ALA:HB3	2.39	0.53
1:B:61:GLN:HE21	1:B:62:VAL:H	1.57	0.53
1:D:222:THR:HB	1:D:237:THR:HG23	1.91	0.53
1:D:34:LEU:HD12	1:D:129:VAL:CG1	2.38	0.53
1:D:227:HIS:HB3	1:D:230:LEU:HB2	1.91	0.53
1:F:161:GLU:CB	1:F:244:HIS:HA	2.38	0.52
1:A:131:THR:HG23	1:A:134:ALA:HB3	1.92	0.52
1:C:62:VAL:HG11	1:C:108:LEU:HA	1.91	0.52
1:D:69:ARG:NH1	1:D:94:TYR:HH	2.07	0.52
1:D:36:THR:HG21	1:D:139:ALA:HB1	1.90	0.52
1:F:165:LEU:HD12	1:F:209:HIS:HB3	1.91	0.52
1:F:191:THR:CA	1:F:207:GLU:O	2.57	0.52
1:A:35:GLU:OE2	1:A:55:ARG:NE	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:VAL:CG2	1:A:124:GLU:O	2.58	0.52
1:C:108:LEU:HD12	1:C:108:LEU:H	1.75	0.52
1:C:118:VAL:HG12	1:C:119:GLN:H	1.75	0.52
1:D:131:THR:O	1:D:132:PHE:HD2	1.93	0.52
1:B:32:GLY:C	1:B:33:GLU:HG2	2.34	0.52
1:F:187:VAL:HG23	1:F:208:PHE:CE2	2.44	0.52
1:B:32:GLY:HA2	1:B:54:TYR:OH	2.11	0.51
1:F:82:LEU:HD21	1:F:112:VAL:CG2	2.37	0.51
1:F:194:ARG:CD	1:F:195:SER:H	2.22	0.51
1:D:34:LEU:HD12	1:D:129:VAL:HG11	1.92	0.51
1:A:159:LEU:O	1:A:242:VAL:HA	2.11	0.51
1:B:52:CYS:HB2	1:B:67:TRP:CZ2	2.45	0.51
1:B:112:VAL:CG1	1:B:113:LEU:N	2.73	0.51
1:C:81:LEU:C	1:C:81:LEU:HD12	2.35	0.51
1:D:141:LEU:HD12	1:D:141:LEU:C	2.36	0.50
1:F:37:SER:HB3	1:F:141:LEU:HD22	1.93	0.50
1:A:60:GLU:HB3	1:A:133:PRO:HD2	1.93	0.50
1:D:151:PRO:HG3	1:D:225:VAL:HG21	1.93	0.50
1:F:151:PRO:CB	1:F:173:ALA:HB2	2.42	0.50
1:C:98:VAL:CG1	1:C:112:VAL:HG11	2.42	0.50
1:D:176:SER:HB2	1:D:201:SER:HA	1.92	0.50
1:D:197:LYS:NZ	1:D:202:ALA:HB2	2.27	0.50
1:E:128:ARG:HA	1:E:138:GLN:HB3	1.94	0.50
1:E:232:GLN:HE21	1:E:232:GLN:CA	2.23	0.50
1:D:131:THR:O	1:D:132:PHE:CD2	2.65	0.50
1:A:70:VAL:HG12	1:A:140:ARG:HH12	1.75	0.49
1:E:66:ALA:HB3	1:E:128:ARG:CZ	2.13	0.49
1:F:103:PRO:HB3	1:F:104:PRO:HA	1.94	0.49
1:F:176:SER:HA	1:F:177:PRO:C	2.37	0.49
1:A:213:SER:O	1:A:242:VAL:HG21	2.12	0.49
1:B:151:PRO:HB2	1:B:236:ILE:HD11	1.94	0.49
1:C:184:ASP:O	1:C:222:THR:HG23	2.13	0.49
1:E:69:ARG:HH22	1:E:121:ASP:HA	1.77	0.49
1:C:43:VAL:HG22	1:C:46:GLN:HG2	1.94	0.49
1:F:65:VAL:HA	1:F:129:VAL:HG23	1.94	0.49
1:B:195:SER:OG	1:B:204:VAL:HG22	2.13	0.49
1:C:227:HIS:CG	1:C:228:PRO:HD2	2.47	0.49
1:E:114:LEU:HD21	1:E:121:ASP:OD1	2.12	0.49
1:E:227:HIS:CG	1:E:228:PRO:HD2	2.47	0.49
1:A:61:GLN:HG3	1:A:62:VAL:N	2.26	0.49
1:D:52:CYS:HB2	1:D:67:TRP:CZ2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:CYS:HB2	1:A:67:TRP:HZ2	1.75	0.49
1:F:105:ARG:HH12	1:F:110:GLY:CA	2.24	0.49
1:C:211:VAL:HG23	1:C:211:VAL:O	2.11	0.49
1:E:69:ARG:NH2	1:E:125:TYR:OH	2.45	0.49
1:C:106:ASN:OD1	1:C:108:LEU:HD12	2.13	0.49
1:A:99:GLU:OE2	1:A:115:ARG:CD	2.61	0.48
1:A:151:PRO:HG3	1:A:225:VAL:HG11	1.95	0.48
1:B:34:LEU:C	1:B:34:LEU:CD2	2.85	0.48
1:B:159:LEU:O	1:B:242:VAL:HA	2.13	0.48
1:C:69:ARG:NH1	1:C:125:TYR:OH	2.46	0.48
1:C:183:TRP:CD1	1:C:191:THR:HG21	2.49	0.48
1:E:51:PRO:HD2	1:E:141:LEU:CD2	2.42	0.48
1:A:211:VAL:HG23	1:A:211:VAL:O	2.11	0.48
1:F:194:ARG:HG2	1:F:194:ARG:NH1	2.27	0.48
1:D:174:GLU:HG3	1:D:203:ALA:HB2	1.95	0.48
1:E:79:LEU:CG	1:E:94:TYR:CZ	2.91	0.48
1:B:60:GLU:HB3	1:B:133:PRO:HD2	1.96	0.48
1:B:188:LYS:HB3	1:B:188:LYS:NZ	2.28	0.48
1:C:51:PRO:HD2	1:C:141:LEU:CD2	2.44	0.48
1:F:153:LEU:HB2	1:F:236:ILE:HG21	1.94	0.48
1:F:194:ARG:HH11	1:F:194:ARG:HG3	1.78	0.48
1:F:222:THR:HB	1:F:237:THR:HB	1.96	0.48
1:A:35:GLU:OE2	1:A:55:ARG:HD2	2.14	0.48
1:C:49:LYS:C	1:C:50:LEU:HD12	2.38	0.48
1:E:60:GLU:CG	1:E:132:PHE:O	2.62	0.48
1:E:151:PRO:HG3	1:E:225:VAL:CG1	2.43	0.48
1:C:112:VAL:HG12	1:C:113:LEU:N	2.29	0.48
1:D:118:VAL:HG12	1:D:119:GLN:H	1.79	0.48
1:F:51:PRO:HA	1:F:111:SER:HA	1.95	0.48
1:A:143:LEU:CD2	1:A:145:VAL:HG23	2.44	0.47
1:E:224:VAL:O	1:E:224:VAL:CG2	2.61	0.47
1:E:239:ILE:HD12	1:E:239:ILE:H	1.78	0.47
1:B:51:PRO:HD2	1:B:141:LEU:HD22	1.95	0.47
1:E:159:LEU:HD21	1:E:167:LEU:O	2.13	0.47
1:A:222:THR:HB	1:A:237:THR:OG1	2.14	0.47
1:E:34:LEU:CD2	1:E:129:VAL:HG11	2.44	0.47
1:F:147:VAL:HG13	1:F:176:SER:H	1.78	0.47
1:B:91:SER:HB2	1:B:92:PRO:HD2	1.97	0.47
1:B:119:GLN:CD	1:B:147:VAL:HG11	2.39	0.47
1:E:69:ARG:HB2	1:E:124:GLU:O	2.15	0.47
1:F:36:THR:HA	1:F:51:PRO:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:40:VAL:CG2	1:F:50:LEU:CD2	2.91	0.47
1:F:48:ALA:HB3	1:F:114:LEU:HB2	1.96	0.47
1:F:52:CYS:HB2	1:F:67:TRP:CZ2	2.50	0.47
1:A:81:LEU:HD12	1:A:81:LEU:C	2.40	0.47
1:F:105:ARG:NH1	1:F:110:GLY:HA3	2.25	0.47
1:A:35:GLU:CD	1:A:55:ARG:HE	2.22	0.46
1:C:52:CYS:HB2	1:C:67:TRP:CZ2	2.49	0.46
1:C:187:VAL:HB	1:C:216:MET:HE2	1.97	0.46
1:E:52:CYS:HB2	1:E:67:TRP:CZ2	2.51	0.46
1:E:230:LEU:CD2	1:E:234:GLN:NE2	2.78	0.46
1:B:83:HIS:NE2	1:C:83:HIS:NE2	2.63	0.46
1:E:114:LEU:HD21	1:E:121:ASP:CG	2.40	0.46
1:F:69:ARG:HB3	1:F:125:TYR:CE1	2.50	0.46
1:A:161:GLU:HG3	1:A:214:ARG:HB2	1.98	0.46
1:F:42:VAL:HG21	1:F:48:ALA:CB	2.45	0.46
1:A:131:THR:HG21	1:A:137:PHE:CE2	2.51	0.46
1:B:224:VAL:O	1:B:224:VAL:HG22	2.16	0.46
1:D:120:ALA:HB2	1:D:200:ARG:NH1	2.21	0.46
1:E:79:LEU:CG	1:E:94:TYR:CE1	2.81	0.46
1:B:231:LEU:CD2	1:B:231:LEU:N	2.78	0.46
1:F:69:ARG:HB3	1:F:125:TYR:CD1	2.51	0.46
1:B:151:PRO:HG2	1:B:234:GLN:HG2	1.96	0.46
1:B:83:HIS:CD2	1:C:83:HIS:HE2	2.34	0.46
1:E:34:LEU:O	1:E:36:THR:HG22	2.15	0.46
1:A:62:VAL:HG13	1:A:108:LEU:HG	1.98	0.45
1:F:233:ASP:OD1	1:F:233:ASP:N	2.46	0.45
1:D:100:GLN:HG3	1:D:101:PRO:HD2	1.97	0.45
1:A:69:ARG:HG3	1:A:125:TYR:CE1	2.51	0.45
1:F:182:THR:OG1	1:F:224:VAL:HG13	2.16	0.45
1:B:82:LEU:HB3	1:B:88:LEU:HD12	1.97	0.45
1:C:103:PRO:HA	1:C:105:ARG:HG3	1.98	0.45
1:E:79:LEU:O	1:E:91:SER:OG	2.35	0.45
1:E:90:VAL:HG21	1:E:94:TYR:C	2.41	0.45
1:E:151:PRO:HB2	1:E:236:ILE:HD11	1.98	0.45
1:C:164:GLY:O	1:C:166:THR:HG23	2.16	0.45
1:A:36:THR:HG21	1:A:139:ALA:HB1	1.98	0.45
1:A:61:GLN:HA	1:A:108:LEU:HD11	1.98	0.45
1:B:53:PHE:HD2	1:B:55:ARG:HH11	1.62	0.45
1:B:147:VAL:O	1:B:147:VAL:HG22	2.17	0.45
1:C:153:LEU:HG	1:C:154:ASN:N	2.29	0.45
1:D:131:THR:HG23	1:D:135:GLY:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:LEU:N	1:B:231:LEU:HD23	2.32	0.45
1:E:42:VAL:HG22	1:E:143:LEU:HD21	1.99	0.45
1:E:195:SER:HA	1:E:204:VAL:HA	1.99	0.45
1:E:36:THR:HG21	1:E:139:ALA:HB1	1.98	0.44
1:F:183:TRP:CZ3	1:F:223:CYS:HB3	2.51	0.44
1:A:156:GLY:HA2	1:A:167:LEU:HD23	2.00	0.44
1:C:69:ARG:HH12	1:C:121:ASP:HA	1.83	0.44
1:A:163:GLN:O	1:A:212:PRO:HD2	2.18	0.44
1:D:118:VAL:CG1	1:D:119:GLN:N	2.80	0.44
1:D:141:LEU:HD12	1:D:141:LEU:O	2.17	0.44
1:D:197:LYS:HZ2	1:D:202:ALA:HB2	1.82	0.44
1:C:131:THR:OG1	1:C:134:ALA:HB3	2.17	0.44
1:C:242:VAL:O	1:C:242:VAL:HG12	2.18	0.44
1:D:98:VAL:CG1	1:D:112:VAL:HG11	2.48	0.44
1:D:119:GLN:O	1:D:119:GLN:HG3	2.18	0.44
1:F:234:GLN:HE21	1:F:234:GLN:N	2.16	0.44
1:D:58:SER:O	1:D:60:GLU:N	2.51	0.44
1:D:174:GLU:CG	1:D:203:ALA:HB2	2.47	0.44
1:E:41:THR:HG23	1:E:144:ARG:CD	2.48	0.44
1:F:68:ALA:HB2	1:F:78:GLU:OE2	2.18	0.44
1:A:116:ASN:O	1:A:118:VAL:HG23	2.17	0.43
1:B:213:SER:O	1:B:216:MET:HG2	2.18	0.43
1:C:98:VAL:HG13	1:C:112:VAL:HG11	2.00	0.43
1:F:65:VAL:HA	1:F:129:VAL:CG2	2.48	0.43
1:E:79:LEU:HG	1:E:94:TYR:CD1	2.49	0.43
1:C:168:ALA:HB2	1:C:210:LEU:HD22	2.00	0.43
1:E:34:LEU:HD21	1:E:129:VAL:CG1	2.46	0.43
1:E:118:VAL:CG1	1:E:119:GLN:N	2.81	0.43
1:C:150:LEU:HD23	1:C:150:LEU:HA	1.76	0.43
1:D:94:TYR:HB3	1:D:98:VAL:CG2	2.48	0.43
1:A:178:ALA:HB2	1:A:202:ALA:HB3	2.00	0.43
1:C:49:LYS:HA	1:C:113:LEU:CD1	2.43	0.43
1:F:235:ARG:HG3	1:F:235:ARG:O	2.18	0.43
1:D:71:ASP:HB2	1:D:75:GLY:O	2.18	0.43
1:F:36:THR:HG21	1:F:139:ALA:HB1	1.99	0.43
1:F:65:VAL:HB	1:F:82:LEU:O	2.18	0.43
1:F:96:GLY:O	1:F:115:ARG:NH2	2.52	0.43
1:F:103:PRO:CB	1:F:104:PRO:CA	2.97	0.43
1:F:151:PRO:HG3	1:F:225:VAL:HG21	2.00	0.43
1:A:118:VAL:HG12	1:A:119:GLN:N	2.33	0.43
1:A:177:PRO:CG	1:A:228:PRO:HG2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:GLN:HB3	1:C:166:THR:HG21	2.01	0.43
1:A:146:LEU:CD2	1:A:177:PRO:HD3	2.48	0.42
1:B:222:THR:HG22	1:B:237:THR:CG2	2.42	0.42
1:E:97:ARG:NH1	1:E:115:ARG:O	2.50	0.42
1:C:49:LYS:HD3	1:C:113:LEU:CD1	2.49	0.42
1:E:143:LEU:HD23	1:E:144:ARG:N	2.35	0.42
1:B:57:ASP:OD1	1:B:57:ASP:C	2.62	0.42
1:B:239:ILE:HD12	1:B:241:HIS:HE2	1.85	0.42
1:D:49:LYS:HD2	1:D:113:LEU:HD11	2.01	0.42
1:D:119:GLN:O	1:D:120:ALA:CB	2.66	0.42
1:C:42:VAL:HG11	1:C:48:ALA:HB2	2.01	0.42
1:C:177:PRO:CG	1:C:228:PRO:HG2	2.49	0.42
1:F:186:GLU:N	1:F:186:GLU:OE2	2.51	0.42
1:A:143:LEU:HD21	1:A:145:VAL:HG23	2.00	0.42
1:D:216:MET:HA	1:D:219:GLN:CG	2.44	0.42
1:E:47:ASP:OD1	1:E:115:ARG:HG3	2.19	0.42
1:F:50:LEU:CD2	1:F:141:LEU:HD11	2.44	0.42
1:F:53:PHE:C	1:F:109:ASP:HB3	2.44	0.42
1:C:49:LYS:HB2	1:C:113:LEU:HD11	2.02	0.42
1:B:106:ASN:OD1	1:B:107:PRO:HD2	2.20	0.42
1:C:53:PHE:CE1	1:C:108:LEU:HB2	2.55	0.42
1:D:36:THR:CG2	1:D:139:ALA:HB1	2.50	0.42
1:D:214:ARG:CZ	1:D:244:HIS:CD2	3.02	0.42
1:E:182:THR:HB	1:E:224:VAL:CG2	2.50	0.42
1:C:35:GLU:OE1	1:C:55:ARG:HG2	2.19	0.42
1:D:65:VAL:HG22	1:D:129:VAL:CG2	2.46	0.42
1:A:85:LYS:HD3	1:A:86:TYR:CZ	2.55	0.42
1:F:234:GLN:HE21	1:F:234:GLN:CA	2.33	0.42
1:A:70:VAL:CG2	1:A:124:GLU:HB3	2.38	0.41
1:A:70:VAL:HG23	1:A:124:GLU:HB2	1.99	0.41
1:A:91:SER:OG	1:A:92:PRO:CD	2.61	0.41
1:A:177:PRO:HD2	1:A:227:HIS:CE1	2.55	0.41
1:F:53:PHE:HE2	1:F:108:LEU:HB3	1.77	0.41
1:A:132:PHE:CD2	1:A:132:PHE:N	2.88	0.41
1:B:151:PRO:HG3	1:B:225:VAL:HG11	2.01	0.41
1:F:91:SER:HB2	1:F:92:PRO:HD2	2.02	0.41
1:C:113:LEU:HD12	1:C:113:LEU:HA	1.74	0.41
1:F:153:LEU:CD2	1:F:236:ILE:HG21	2.49	0.41
1:F:159:LEU:HD23	1:F:166:THR:HB	2.02	0.41
1:A:131:THR:HG21	1:A:137:PHE:HE2	1.86	0.41
1:A:194:ARG:HA	1:A:194:ARG:HD3	1.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:GLU:OE2	1:A:81:LEU:HD23	2.20	0.41
1:C:101:PRO:O	1:C:105:ARG:CD	2.51	0.41
1:E:118:VAL:HG12	1:E:119:GLN:H	1.83	0.41
1:F:153:LEU:HD23	1:F:236:ILE:HG21	1.96	0.41
1:F:222:THR:HB	1:F:237:THR:CB	2.50	0.41
1:D:94:TYR:HB3	1:D:98:VAL:HG23	2.03	0.41
1:F:152:SER:N	1:F:172:THR:O	2.52	0.41
1:F:165:LEU:HA	1:F:210:LEU:O	2.21	0.41
1:C:178:ALA:HB2	1:C:202:ALA:HB3	2.02	0.41
1:D:213:SER:O	1:D:242:VAL:HG11	2.21	0.41
1:E:149:PRO:HG2	1:E:234:GLN:HE21	1.84	0.41
1:F:109:ASP:OD2	1:F:109:ASP:O	2.39	0.41
1:F:184:ASP:O	1:F:185:THR:OG1	2.28	0.41
1:A:35:GLU:OE2	1:A:55:ARG:CD	2.69	0.41
1:F:68:ALA:HB2	1:F:78:GLU:CD	2.46	0.41
1:A:227:HIS:CD2	1:A:228:PRO:HD2	2.56	0.40
1:B:116:ASN:N	1:B:116:ASN:HD22	2.18	0.40
1:B:73:GLY:C	1:B:75:GLY:H	2.28	0.40
1:C:118:VAL:CG1	1:C:119:GLN:N	2.85	0.40
1:D:194:ARG:NH2	1:E:150:LEU:HG	2.36	0.40
1:D:224:VAL:O	1:D:224:VAL:CG2	2.70	0.40
1:F:210:LEU:C	1:F:210:LEU:HD12	2.46	0.40
1:A:183:TRP:CD1	1:A:191:THR:HG21	2.57	0.40
1:A:122:GLU:OE2	1:A:144:ARG:NH1	2.43	0.40
1:B:34:LEU:HD21	1:B:52:CYS:SG	2.61	0.40
1:C:70:VAL:HG21	1:C:124:GLU:HB3	2.01	0.40
1:E:60:GLU:HG2	1:E:133:PRO:HD2	2.04	0.40
1:E:184:ASP:O	1:E:222:THR:HG23	2.21	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	213/218 (98%)	206 (97%)	7 (3%)	0	100	100
1	B	211/218 (97%)	204 (97%)	7 (3%)	0	100	100
1	C	209/218 (96%)	199 (95%)	10 (5%)	0	100	100
1	D	210/218 (96%)	197 (94%)	12 (6%)	1 (0%)	24	57
1	E	206/218 (94%)	190 (92%)	16 (8%)	0	100	100
1	F	181/218 (83%)	167 (92%)	14 (8%)	0	100	100
All	All	1230/1308 (94%)	1163 (95%)	66 (5%)	1 (0%)	48	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	59	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	176/181 (97%)	156 (89%)	20 (11%)	5	24
1	B	176/181 (97%)	144 (82%)	32 (18%)	2	10
1	C	171/181 (94%)	144 (84%)	27 (16%)	2	15
1	D	169/181 (93%)	137 (81%)	32 (19%)	1	9
1	E	165/181 (91%)	139 (84%)	26 (16%)	2	15
1	F	150/181 (83%)	119 (79%)	31 (21%)	1	6
All	All	1007/1086 (93%)	839 (83%)	168 (17%)	2	13

All (168) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	ASP
1	A	100	GLN
1	A	114	LEU
1	A	131	THR

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Mol	Chain	Res	Type
1	A	147	VAL
1	A	152	SER
1	A	161	GLU
1	A	192	SER
1	A	197	LYS
1	A	200	ARG
1	A	204	VAL
1	A	205	THR
1	A	210	LEU
1	A	222	THR
1	A	224	VAL
1	A	225	VAL
1	A	236	ILE
1	A	242	VAL
1	A	243	SER
1	A	244	HIS
1	B	33	GLU
1	B	34	LEU
1	B	41	THR
1	B	43	VAL
1	B	58	SER
1	B	61	GLN
1	B	71	ASP
1	B	81	LEU
1	B	98	VAL
1	B	99	GLU
1	B	100	GLN
1	B	105	ARG
1	B	114	LEU
1	B	116	ASN
1	B	129	VAL
1	B	131	THR
1	B	140	ARG
1	B	144	ARG
1	B	147	VAL
1	B	152	SER
1	B	161	GLU
1	B	188	LYS
1	B	192	SER
1	B	197	LYS
1	B	200	ARG
1	B	205	THR

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Mol	Chain	Res	Type
1	B	222	THR
1	B	224	VAL
1	B	225	VAL
1	B	231	LEU
1	B	236	ILE
1	B	244	HIS
1	C	35	GLU
1	C	49	LYS
1	C	50	LEU
1	C	62	VAL
1	C	70	VAL
1	C	95	GLU
1	C	97	ARG
1	C	108	LEU
1	C	113	LEU
1	C	129	VAL
1	C	130	SER
1	C	131	THR
1	C	147	VAL
1	C	152	SER
1	C	153	LEU
1	C	192	SER
1	C	193	SER
1	C	199	SER
1	C	205	THR
1	C	210	LEU
1	C	213	SER
1	C	217	ASN
1	C	222	THR
1	C	224	VAL
1	C	225	VAL
1	C	230	LEU
1	C	237	THR
1	D	38	ASP
1	D	42	VAL
1	D	43	VAL
1	D	46	GLN
1	D	58	SER
1	D	69	ARG
1	D	70	VAL
1	D	78	GLU
1	D	79	LEU

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Mol	Chain	Res	Type
1	D	82	LEU
1	D	114	LEU
1	D	124	GLU
1	D	131	THR
1	D	138	GLN
1	D	142	ARG
1	D	143	LEU
1	D	146	LEU
1	D	147	VAL
1	D	154	ASN
1	D	161	GLU
1	D	172	THR
1	D	190	THR
1	D	197	LYS
1	D	210	LEU
1	D	214	ARG
1	D	215	SER
1	D	222	THR
1	D	224	VAL
1	D	225	VAL
1	D	231	LEU
1	D	236	ILE
1	D	239	ILE
1	E	41	THR
1	E	42	VAL
1	E	43	VAL
1	E	60	GLU
1	E	62	VAL
1	E	70	VAL
1	E	74	GLU
1	E	82	LEU
1	E	100	GLN
1	E	114	LEU
1	E	129	VAL
1	E	138	GLN
1	E	143	LEU
1	E	147	VAL
1	E	152	SER
1	E	182	THR
1	E	197	LYS
1	E	201	SER
1	E	204	VAL

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Mol	Chain	Res	Type
1	E	205	THR
1	E	210	LEU
1	E	222	THR
1	E	225	VAL
1	E	230	LEU
1	E	236	ILE
1	E	239	ILE
1	F	44	LEU
1	F	70	VAL
1	F	99	GLU
1	F	109	ASP
1	F	111	SER
1	F	112	VAL
1	F	113	LEU
1	F	114	LEU
1	F	118	VAL
1	F	129	VAL
1	F	131	THR
1	F	140	ARG
1	F	147	VAL
1	F	152	SER
1	F	153	LEU
1	F	159	LEU
1	F	160	GLU
1	F	166	THR
1	F	170	SER
1	F	171	CYS
1	F	184	ASP
1	F	191	THR
1	F	194	ARG
1	F	204	VAL
1	F	205	THR
1	F	214	ARG
1	F	222	THR
1	F	224	VAL
1	F	225	VAL
1	F	230	LEU
1	F	234	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	219	GLN
1	B	61	GLN
1	B	116	ASN
1	B	217	ASN
1	B	219	GLN
1	C	64	GLN
1	C	89	HIS
1	C	119	GLN
1	C	209	HIS
1	C	217	ASN
1	C	219	GLN
1	C	241	HIS
1	D	46	GLN
1	D	89	HIS
1	D	219	GLN
1	D	241	HIS
1	D	244	HIS
1	E	46	GLN
1	E	61	GLN
1	E	64	GLN
1	E	198	HIS
1	E	217	ASN
1	E	232	GLN
1	E	234	GLN
1	F	100	GLN
1	F	119	GLN
1	F	198	HIS
1	F	234	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	215/218 (98%)	-0.36	0	100 100	35, 59, 99, 131	0
1	B	213/218 (97%)	-0.39	0	100 100	40, 58, 89, 133	0
1	C	211/218 (96%)	-0.30	0	100 100	39, 58, 121, 148	0
1	D	212/218 (97%)	-0.23	0	100 100	68, 107, 148, 189	0
1	E	208/218 (95%)	0.02	4 (1%)	66 41	55, 118, 206, 260	0
1	F	189/218 (86%)	0.78	13 (6%)	23 14	114, 145, 192, 263	0
All	All	1248/1308 (95%)	-0.10	17 (1%)	73 48	35, 84, 168, 263	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	107	PRO	4.9
1	F	142	ARG	3.8
1	E	87	GLY	3.5
1	F	169	ALA	3.5
1	F	225	VAL	3.2
1	F	178	ALA	3.0
1	F	64	GLN	2.9
1	F	228	PRO	2.8
1	F	40	VAL	2.7
1	F	181	VAL	2.6
1	E	110	GLY	2.6
1	F	179	PRO	2.5
1	F	158	ALA	2.5
1	F	173	ALA	2.1
1	E	86	TYR	2.1
1	F	93	ALA	2.1
1	E	82	LEU	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.