



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:33 AM UTC

PDB ID : 3QDZ / pdb_00003qdz
Title : Crystal structure of the human thrombin mutant D102N in complex with the extracellular fragment of human PAR4.
Authors : Gandhi, P.; Chen, Z.; Appelbaum, E.; Zapata, F.; Di Cera, E.
Deposited on : 2011-01-19
Resolution : 2.80 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

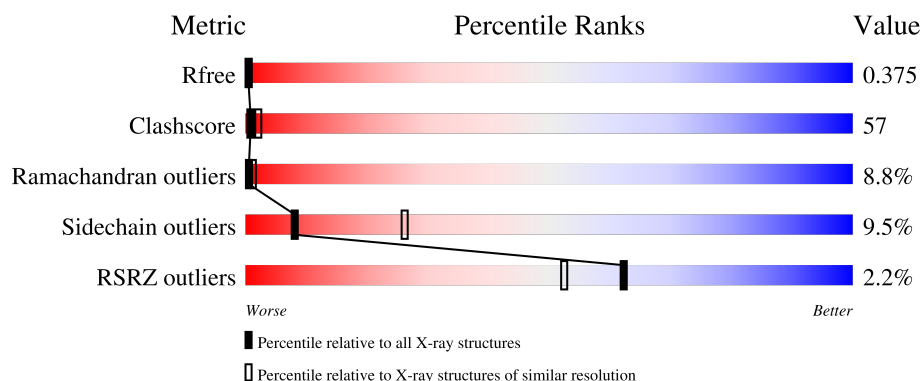
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

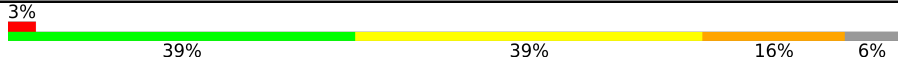
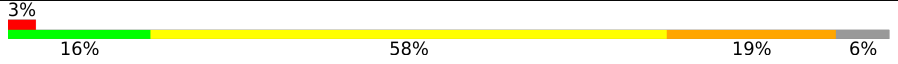
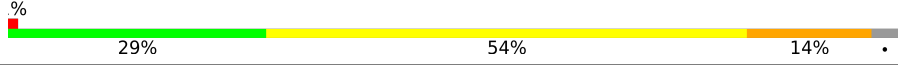
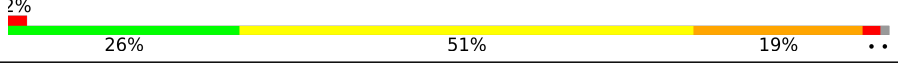
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	31	
1	C	31	
2	B	259	
2	D	259	
3	E	9	

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Mol	Chain	Length	Quality of chain
3	F	9	22%22%56%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4677 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 Thrombin light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	29	Total	C	N	O	S	0	0	0
			239	149	38	51	1			
1	C	29	Total	C	N	O	S	0	0	0
			239	149	38	51	1			

- Molecule 2 is a protein called Thrombin heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	251	Total	C	N	O	S	0	0	0
			2028	1295	358	361	14			
2	D	257	Total	C	N	O	S	0	0	0
			2073	1324	366	369	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	102	ASN	ASP	engineered mutation	UNP P00734
D	102	ASN	ASP	engineered mutation	UNP P00734

- Molecule 3 is a protein called Proteinase-activated receptor 4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	9	Total	C	N	O	0	0	0
			67	43	12	12			
3	F	4	Total	C	N	O	0	0	0
			31	19	7	5			

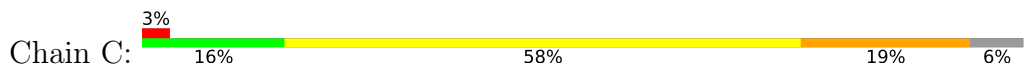
3 Residue-property plots

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

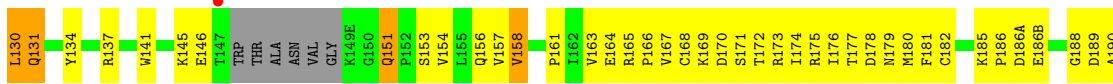
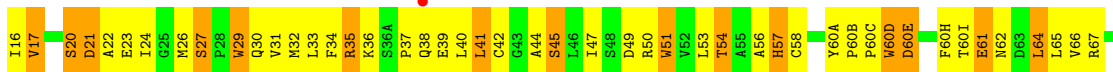
- Molecule 1: Thrombin light chain



- Molecule 1: Thrombin light chain

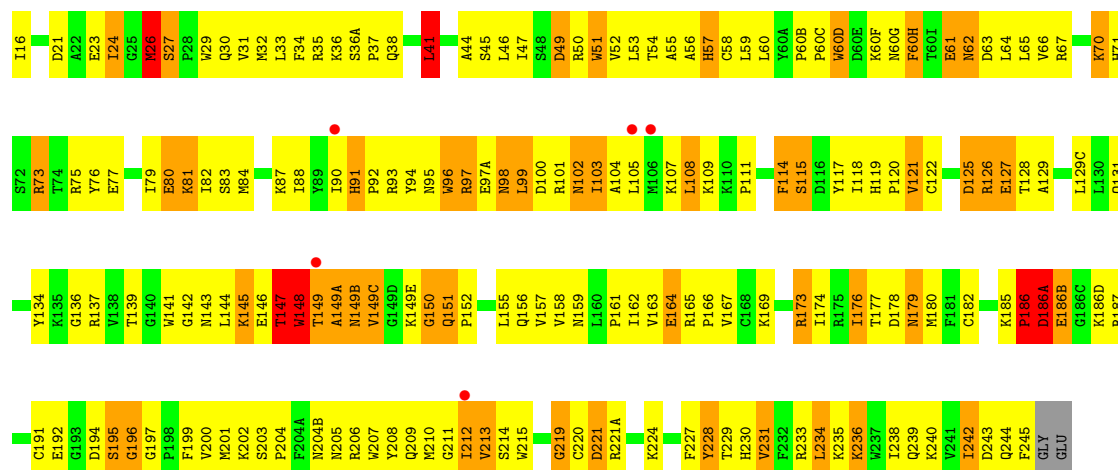


- Molecule 2: Thrombin heavy chain



- Molecule 2: Thrombin heavy chain





● Molecule 3: Proteinase-activated receptor 4



● Molecule 3: Proteinase-activated receptor 4



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	91.41Å 102.88Å 146.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.80 40.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	86.0 (40.00-2.80) 86.3 (40.00-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.84 (at 2.81Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.304 , 0.364 0.315 , 0.375	Depositor DCC
R_{free} test set	796 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	55.4	Xtriage
Anisotropy	1.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 75.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4677	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% 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	A	0.54	0/241	1.11	3/321 (0.9%)
1	C	0.43	0/241	0.97	2/321 (0.6%)
2	B	0.62	0/2080	1.10	17/2810 (0.6%)
2	D	0.58	1/2128 (0.0%)	1.16	23/2879 (0.8%)
3	E	0.67	0/69	1.36	0/94
3	F	0.68	0/32	1.17	0/41
All	All	0.59	1/4791 (0.0%)	1.12	45/6466 (0.7%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	148	TRP	CA-C	6.72	1.61	1.52

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	62	ASN	N-CA-C	8.30	121.44	111.82
2	D	96	TRP	N-CA-C	-7.73	103.46	113.12
2	D	149	THR	N-CA-C	7.69	127.19	110.80
2	B	110	LYS	CA-C-N	7.57	129.31	119.84
2	B	110	LYS	C-N-CA	7.57	129.31	119.84
2	D	61	GLU	N-CA-C	7.15	119.74	111.02
2	D	91	HIS	CA-C-N	6.74	126.25	119.24
2	D	91	HIS	C-N-CA	6.74	126.25	119.24
2	D	179	ASN	N-CA-C	-6.56	105.31	113.38
2	B	22	ALA	N-CA-C	6.53	118.98	110.43
2	D	186(B)	GLU	N-CA-C	-6.40	103.78	113.89
2	D	219	GLY	N-CA-C	-6.39	103.61	112.18
2	B	60(D)	TRP	N-CA-C	-6.33	102.95	111.87
2	D	147	THR	N-CA-C	6.26	124.13	110.80
2	B	64	LEU	N-CA-C	6.13	118.73	109.41
2	D	49	ASP	N-CA-C	-6.04	104.99	112.90
2	D	51	TRP	N-CA-C	5.93	118.30	108.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	35	ARG	N-CA-C	-5.76	100.61	109.76
2	B	27	SER	CA-C-N	5.70	126.97	119.84
2	B	27	SER	C-N-CA	5.70	126.97	119.84
2	D	26	MET	N-CA-C	-5.68	106.52	113.50
2	B	97(A)	GLU	N-CA-C	5.66	120.19	112.04
1	A	14(J)	TYR	N-CA-C	-5.59	106.13	113.12
1	A	14	ASP	N-CA-C	-5.51	103.42	110.53
2	D	242	ILE	N-CA-C	-5.46	106.50	113.22
2	B	190	ALA	N-CA-C	-5.45	101.91	110.36
2	B	29	TRP	N-CA-C	-5.42	105.97	112.59
2	B	204	PRO	N-CA-C	-5.41	107.71	114.20
2	B	62	ASN	N-CA-C	5.41	119.63	113.19
1	A	14(K)	ILE	N-CA-C	-5.38	105.60	110.82
2	D	196	GLY	N-CA-C	-5.38	108.21	115.36
1	C	4	ARG	CA-C-N	5.37	126.55	119.84
1	C	4	ARG	C-N-CA	5.37	126.55	119.84
2	D	60(D)	TRP	N-CA-C	-5.29	107.00	112.93
2	D	145	LYS	N-CA-C	-5.29	101.66	109.18
2	B	17	VAL	N-CA-C	5.28	116.06	108.93
2	B	51	TRP	N-CA-C	5.28	118.16	109.72
2	D	149	THR	CB-CA-C	-5.25	99.98	110.42
2	D	71	HIS	N-CA-C	-5.25	105.09	112.12
2	B	21	ASP	N-CA-C	-5.21	100.91	109.40
2	B	60(H)	PHE	N-CA-C	5.20	117.39	110.53
2	D	41	LEU	N-CA-C	5.17	119.48	112.04
2	D	73	ARG	N-CA-C	-5.10	106.74	113.12
2	D	27	SER	CA-C-N	5.03	124.75	119.82
2	D	27	SER	C-N-CA	5.03	124.75	119.82

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	A	239	0	235	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	239	0	235	41	0
2	B	2028	0	2005	203	0
2	D	2073	0	2046	285	0
3	E	67	0	72	17	0
3	F	31	0	32	9	0
All	All	4677	0	4625	527	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 57.

All (527) 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)
3:E:39:THR:OG1	3:E:40:PRO:HD3	1.26	1.28
2:D:173:ARG:HB3	2:D:173:ARG:HH11	1.13	1.13
3:E:39:THR:HG1	3:E:40:PRO:HD3	1.23	1.00
2:D:55:ALA:HB1	2:D:102:ASN:HD21	1.22	0.99
2:D:173:ARG:HB3	2:D:173:ARG:NH1	1.78	0.98
2:D:67:ARG:HG2	2:D:82:ILE:HG12	1.48	0.96
2:D:195:SER:HA	2:D:213:VAL:HB	1.47	0.95
2:D:31:VAL:HB	2:D:44:ALA:HB3	1.50	0.94
2:B:157:VAL:HG22	2:B:158:VAL:H	1.34	0.92
2:B:95:ASN:HD21	2:B:97(A):GLU:HB3	1.35	0.91
2:B:129(B):SER:HB2	2:D:149:THR:HG22	1.54	0.88
2:D:235:LYS:HA	2:D:238:ILE:HD12	1.53	0.88
2:D:136:GLY:HA3	2:D:199:PHE:CE1	2.09	0.87
2:B:244:GLN:O	2:B:245:PHE:HB2	1.73	0.86
2:D:90:ILE:HG22	2:D:91:HIS:H	1.42	0.85
2:B:128:THR:HG22	2:B:129:ALA:N	1.90	0.85
2:B:73:ARG:HH12	2:B:151:GLN:HG3	1.42	0.85
2:D:238:ILE:HG22	2:D:242:ILE:HD11	1.59	0.83
3:E:39:THR:OG1	3:E:40:PRO:CD	2.22	0.82
2:D:91:HIS:HE1	2:D:93:ARG:HB2	1.44	0.82
2:D:129:ALA:HA	2:D:210:MET:HE1	1.61	0.81
2:D:32:MET:HG2	2:D:141:TRP:CZ3	2.16	0.81
2:B:101:ARG:O	2:B:103:ILE:HG22	1.81	0.80
2:D:81:LYS:HD2	2:D:81:LYS:N	1.97	0.79
2:D:235:LYS:HA	2:D:238:ILE:CD1	2.12	0.79
1:A:4:ARG:HB2	1:A:8:GLU:OE1	1.83	0.78
2:D:169:LYS:HG3	2:D:176:ILE:HD11	1.63	0.78
2:D:50:ARG:HD2	2:D:111:PRO:HB3	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:157:VAL:HG22	2:B:158:VAL:N	1.96	0.78
2:B:17:VAL:HG11	2:B:221:ASP:HB2	1.66	0.77
1:C:14(D):ARG:HD3	1:C:14(H):GLU:HG3	1.66	0.77
2:D:21:ASP:HA	2:D:156:GLN:HE22	1.49	0.76
2:B:33:LEU:HD11	2:B:64:LEU:HD13	1.67	0.76
2:D:99:LEU:HD21	3:F:46:PRO:HB3	1.67	0.76
2:D:50:ARG:HH21	2:D:107:LYS:HE3	1.49	0.76
2:B:86:GLU:O	2:B:87:LYS:HB2	1.85	0.76
2:D:94:TYR:HD1	2:D:100:ASP:O	1.68	0.76
3:E:41:SER:O	3:E:42:ILE:HG13	1.86	0.76
2:D:149(A):ALA:C	2:D:149(C):VAL:H	1.91	0.76
1:C:14:ASP:CG	2:D:26:MET:HE3	2.11	0.76
2:D:236:LYS:NZ	2:D:236:LYS:HA	2.00	0.75
2:B:99:LEU:HD21	3:E:46:PRO:HB3	1.68	0.75
2:D:61:GLU:OE2	2:D:87:LYS:HD2	1.85	0.75
2:D:90:ILE:HG22	2:D:91:HIS:N	2.02	0.74
2:D:195:SER:OG	3:F:47:ARG:C	2.31	0.74
2:D:47:ILE:HD11	2:D:51:TRP:HB3	1.70	0.73
2:B:60(D):TRP:CH2	3:E:46:PRO:HG2	2.23	0.73
2:D:117:TYR:C	2:D:118:ILE:HD12	2.13	0.73
2:D:45:SER:CB	2:D:209:GLN:HE22	2.01	0.73
2:B:199:PHE:HD1	2:B:199:PHE:O	1.73	0.72
2:D:186(A):ASP:OD1	2:D:186(A):ASP:O	2.07	0.72
2:D:234:LEU:O	2:D:238:ILE:HG13	1.90	0.72
2:D:65:LEU:HD12	2:D:84:MET:HE1	1.72	0.71
2:B:204(A):PHE:HZ	2:D:149:THR:H	1.36	0.71
2:B:53:LEU:O	2:B:54:THR:HB	1.91	0.71
1:A:10:LYS:O	1:A:12:LEU:HD13	1.91	0.70
2:D:60(H):PHE:CD1	2:D:64:LEU:HD21	2.27	0.70
2:D:49:ASP:OD2	2:D:50:ARG:HG2	1.91	0.69
2:D:31:VAL:HB	2:D:44:ALA:CB	2.22	0.69
2:B:61:GLU:HG2	2:B:87:LYS:HA	1.72	0.69
2:B:99:LEU:HG	2:B:102:ASN:ND2	2.08	0.69
2:D:145:LYS:HE3	2:D:149(E):LYS:HD2	1.73	0.69
2:B:56:ALA:HB2	2:B:103:ILE:N	2.08	0.69
2:B:129(B):SER:HB2	2:D:149:THR:CG2	2.21	0.69
2:B:186:PRO:O	2:B:186(A):ASP:HB2	1.91	0.69
2:D:23:GLU:HB2	2:D:26:MET:HG3	1.72	0.69
2:D:55:ALA:HB1	2:D:102:ASN:ND2	2.04	0.68
2:B:157:VAL:CG2	2:B:158:VAL:H	2.05	0.68
1:A:7:PHE:HD1	1:A:12:LEU:O	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:29:TRP:CD2	2:D:121:VAL:HB	2.29	0.68
2:D:65:LEU:HD21	2:D:82:ILE:HG23	1.75	0.68
2:B:92:PRO:HG2	2:B:93:ARG:H	1.60	0.67
2:D:149(A):ALA:O	2:D:149(C):VAL:N	2.28	0.67
2:D:98:ASN:OD1	2:D:100:ASP:HB2	1.94	0.67
2:B:178:ASP:O	2:B:233:ARG:HG3	1.94	0.66
2:D:195:SER:HB3	3:F:47:ARG:HB2	1.77	0.66
2:B:129(C):LEU:O	2:B:131:GLN:N	2.28	0.66
2:B:130:LEU:HD22	2:B:230:HIS:NE2	2.10	0.65
2:D:91:HIS:CE1	2:D:93:ARG:HB2	2.29	0.65
2:B:195:SER:HB3	3:E:47:ARG:HB2	1.78	0.65
2:D:83:SER:O	2:D:84:MET:HE2	1.96	0.65
2:D:149(B):ASN:C	2:D:149(C):VAL:HG22	2.21	0.65
2:B:71:HIS:CD2	2:B:154:VAL:HG13	2.32	0.65
2:B:217:GLU:HB3	3:E:42:ILE:CG2	2.27	0.64
2:D:96:TRP:H	2:D:96:TRP:CD1	2.12	0.64
2:D:126:ARG:CZ	2:D:126:ARG:HB2	2.26	0.64
2:D:211:GLY:HA2	2:D:229:THR:O	1.97	0.64
2:B:204(B):ASN:OD1	2:B:206:ARG:HG3	1.96	0.64
2:D:57:HIS:ND1	3:F:46:PRO:HB2	2.13	0.64
3:E:43:LEU:HD23	3:E:44:PRO:HD2	1.78	0.64
2:D:34:PHE:HE1	2:D:38:GLN:HE21	1.44	0.64
1:A:3:LEU:HD21	2:B:206:ARG:HG2	1.79	0.64
2:B:34:PHE:HE2	2:B:67:ARG:HG3	1.63	0.64
2:D:91:HIS:CE1	2:D:93:ARG:H	2.15	0.64
2:D:200:VAL:HA	2:D:209:GLN:HA	1.79	0.64
2:B:34:PHE:HD1	2:B:39:GLU:O	1.79	0.64
2:B:47:ILE:HD11	2:B:51:TRP:HB2	1.80	0.64
2:D:149(A):ALA:C	2:D:149(C):VAL:N	2.55	0.63
2:D:67:ARG:HD2	2:D:80:GLU:OE1	1.98	0.63
2:B:56:ALA:HB2	2:B:103:ILE:H	1.63	0.63
2:D:149(B):ASN:O	2:D:149(C):VAL:HG22	1.98	0.63
2:B:34:PHE:CE2	2:B:67:ARG:HG3	2.33	0.63
2:D:204(B):ASN:HD22	2:D:204(B):ASN:H	1.45	0.63
2:B:237:TRP:O	2:B:241:VAL:HG23	1.98	0.62
2:D:99:LEU:N	2:D:99:LEU:HD12	2.14	0.62
2:B:127:GLU:HG2	2:D:148:TRP:CZ2	2.35	0.62
1:C:12:LEU:HD13	1:C:13:GLU:N	2.14	0.62
2:D:179:ASN:O	2:D:230:HIS:N	2.22	0.62
2:D:57:HIS:HD2	2:D:58:CYS:N	1.97	0.62
2:D:144:LEU:HB2	2:D:149(E):LYS:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:86:GLU:HB2	2:B:109:LYS:HA	1.82	0.61
2:D:139:THR:HG22	2:D:157:VAL:HB	1.82	0.61
2:D:67:ARG:HD2	2:D:70:LYS:HD2	1.81	0.61
2:D:147:THR:C	2:D:148:TRP:CD1	2.79	0.61
2:D:129(C):LEU:HD22	2:D:201:MET:HE2	1.82	0.61
2:B:199:PHE:O	2:B:199:PHE:CD1	2.53	0.61
2:D:114:PHE:O	2:D:115:SER:HB3	1.99	0.61
2:B:236:LYS:HB3	2:B:236:LYS:HZ2	1.65	0.61
2:D:146:GLU:HG3	2:D:220:CYS:H	1.66	0.61
2:D:50:ARG:CD	2:D:111:PRO:HB3	2.29	0.60
2:B:99:LEU:HG	2:B:102:ASN:HD22	1.66	0.60
2:D:128:THR:O	2:D:129(C):LEU:HB2	2.00	0.60
1:A:14:ASP:HB2	2:B:26:MET:HE3	1.84	0.60
2:B:215:TRP:HA	3:E:47:ARG:HG3	1.83	0.60
2:D:30:GLN:HG2	2:D:155:LEU:HD22	1.83	0.60
2:D:65:LEU:HD23	2:D:66:VAL:N	2.16	0.60
2:D:84:MET:HB3	2:D:109:LYS:HZ2	1.66	0.60
2:D:99:LEU:O	2:D:102:ASN:HB2	2.01	0.60
2:D:233:ARG:C	2:D:234:LEU:HD23	2.26	0.60
2:B:177:THR:OG1	2:B:180:MET:HE3	2.01	0.60
2:D:236:LYS:HA	2:D:236:LYS:HZ2	1.66	0.60
2:D:244:GLN:O	2:D:245:PHE:C	2.45	0.60
2:D:147:THR:C	2:D:148:TRP:HD1	2.09	0.60
2:D:83:SER:C	2:D:84:MET:HE2	2.27	0.59
2:D:186:PRO:O	2:D:186(A):ASP:HB3	2.03	0.59
2:B:49:ASP:HB3	2:B:114:PHE:CZ	2.37	0.59
2:D:53:LEU:HD12	2:D:54:THR:H	1.67	0.59
1:C:7:PHE:HB3	1:C:12:LEU:O	2.01	0.59
2:D:67:ARG:NH2	2:D:80:GLU:OE1	2.30	0.59
2:B:35:ARG:HA	2:B:64:LEU:HD23	1.83	0.59
2:D:227:PHE:C	2:D:228:TYR:CD1	2.81	0.59
2:D:236:LYS:HA	2:D:236:LYS:HZ3	1.66	0.59
2:B:238:ILE:O	2:B:242:ILE:HG13	2.03	0.59
2:D:30:GLN:HG2	2:D:155:LEU:CD2	2.33	0.59
2:D:147:THR:O	2:D:148:TRP:CB	2.51	0.59
2:B:50:ARG:HG2	2:B:111:PRO:HB3	1.85	0.59
2:D:50:ARG:NH1	2:D:111:PRO:HD3	2.18	0.59
2:D:204(B):ASN:H	2:D:204(B):ASN:ND2	2.01	0.59
2:D:50:ARG:HH11	2:D:111:PRO:HD3	1.68	0.58
2:B:60(B):PRO:HB2	2:B:60(C):PRO:HD3	1.85	0.58
1:C:1:CYS:HB2	2:D:122:CYS:N	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:239:GLN:O	2:D:243:ASP:N	2.36	0.58
2:B:165:ARG:N	2:B:166:PRO:HD2	2.18	0.58
2:D:147:THR:O	2:D:148:TRP:CD1	2.56	0.58
2:B:177:THR:H	2:B:180:MET:HE3	1.68	0.58
2:D:77:GLU:HB3	2:D:79:ILE:HB	1.85	0.58
2:B:157:VAL:CG2	2:B:158:VAL:N	2.66	0.58
2:D:195:SER:HG	3:F:47:ARG:C	2.10	0.58
2:D:147:THR:O	2:D:148:TRP:HB2	2.03	0.58
2:D:239:GLN:HA	2:D:242:ILE:HD12	1.86	0.58
2:B:51:TRP:NE1	2:B:242:ILE:HG23	2.19	0.58
2:D:90:ILE:CG2	2:D:91:HIS:H	2.13	0.58
2:D:53:LEU:HD12	2:D:104:ALA:O	2.04	0.57
1:A:3:LEU:CD2	2:B:206:ARG:HG2	2.35	0.57
2:B:67:ARG:HD3	2:B:82:ILE:HG12	1.86	0.57
2:D:228:TYR:CD1	2:D:228:TYR:N	2.72	0.57
2:D:21:ASP:HA	2:D:156:GLN:NE2	2.18	0.57
2:D:59:LEU:O	2:D:60(F):LYS:HG2	2.05	0.57
2:D:95:ASN:OD1	2:D:97:ARG:HB2	2.04	0.57
2:B:203:SER:HB3	2:B:204(B):ASN:ND2	2.20	0.57
2:D:60(H):PHE:HB3	2:D:64:LEU:HD11	1.87	0.57
2:D:16:ILE:HD11	2:D:139:THR:C	2.30	0.56
2:B:77(A):ARG:O	2:B:78:ASN:HB2	2.05	0.56
2:B:228:TYR:CD1	2:B:228:TYR:N	2.73	0.56
2:D:45:SER:O	2:D:52:VAL:HG13	2.05	0.56
2:B:98:ASN:OD1	2:B:100:ASP:HB2	2.06	0.56
2:B:170:ASP:C	2:B:172:THR:H	2.14	0.56
2:D:55:ALA:CB	2:D:102:ASN:HD21	2.08	0.56
2:B:75:ARG:CZ	2:B:77:GLU:HG3	2.35	0.56
3:E:42:ILE:O	3:E:42:ILE:HG22	2.05	0.56
2:D:57:HIS:HA	2:D:60:LEU:O	2.05	0.56
2:B:56:ALA:HB3	2:B:102:ASN:HA	1.88	0.56
2:B:73:ARG:HH22	2:B:151:GLN:HG2	1.70	0.55
2:D:98:ASN:O	2:D:99:LEU:C	2.49	0.55
2:D:169:LYS:HA	2:D:176:ILE:CD1	2.37	0.55
2:B:26:MET:HE2	2:B:137:ARG:NH1	2.22	0.55
2:B:35:ARG:N	2:B:41:LEU:HD11	2.22	0.55
2:B:125:ASP:O	2:B:126:ARG:C	2.48	0.55
2:D:178:ASP:O	2:D:233:ARG:HG3	2.06	0.55
2:B:157:VAL:O	2:B:158:VAL:HB	2.06	0.55
2:B:17:VAL:HG11	2:B:221:ASP:CB	2.35	0.55
2:D:73:ARG:HH12	2:D:151:GLN:HE21	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:29:TRP:CG	2:B:121:VAL:HB	2.42	0.54
2:B:50:ARG:NH1	2:B:107:LYS:HE3	2.22	0.54
2:B:60(I):THR:O	2:B:61:GLU:C	2.51	0.54
2:B:91:HIS:CE1	2:B:101:ARG:HD3	2.42	0.54
1:C:14(D):ARG:HD3	1:C:14(D):ARG:O	2.07	0.54
2:D:50:ARG:NH2	2:D:107:LYS:HE3	2.18	0.54
2:D:169:LYS:CG	2:D:176:ILE:HD11	2.37	0.54
2:B:175:ARG:HH12	2:B:177:THR:CG2	2.20	0.54
2:D:75:ARG:HG2	2:D:76:TYR:H	1.72	0.54
2:D:95:ASN:HD21	2:D:97(A):GLU:HB3	1.73	0.54
2:D:75:ARG:HG2	2:D:76:TYR:N	2.22	0.54
2:B:31:VAL:HB	2:B:44:ALA:HB3	1.90	0.53
2:B:176:ILE:HG21	2:B:227:PHE:CE2	2.43	0.53
2:D:67:ARG:CD	2:D:80:GLU:OE1	2.57	0.53
2:B:34:PHE:HB2	2:B:65:LEU:HD23	1.89	0.53
2:B:95:ASN:ND2	2:B:97(A):GLU:HB3	2.15	0.53
2:D:206:ARG:HB2	2:D:208:TYR:HE2	1.72	0.53
1:C:12:LEU:HD13	1:C:13:GLU:H	1.73	0.53
2:B:60(A):TYR:CE2	2:B:60(C):PRO:HB2	2.43	0.53
2:D:177:THR:OG1	2:D:179:ASN:ND2	2.41	0.53
2:D:60(G):ASN:OD1	2:D:60(G):ASN:O	2.26	0.53
2:B:199:PHE:CD1	2:B:199:PHE:C	2.87	0.53
1:C:8:GLU:HG3	2:D:205:ASN:HD21	1.73	0.53
2:B:16:ILE:HD12	2:B:156:GLN:O	2.08	0.53
2:B:204(B):ASN:HD22	2:B:204(B):ASN:H	1.55	0.53
2:B:211:GLY:HA2	2:B:229:THR:O	2.08	0.53
1:C:14(G):LEU:HD23	1:C:14(J):TYR:HE2	1.74	0.52
2:D:125:ASP:O	2:D:127:GLU:N	2.42	0.52
1:C:4:ARG:HB2	1:C:8:GLU:OE1	2.09	0.52
1:C:7:PHE:CE2	1:C:14:ASP:HB3	2.44	0.52
2:B:215:TRP:HB3	3:E:46:PRO:HA	1.90	0.52
2:B:227:PHE:C	2:B:228:TYR:CD1	2.87	0.52
2:D:44:ALA:HA	2:D:196:GLY:O	2.09	0.52
2:D:94:TYR:CG	2:D:94:TYR:O	2.63	0.52
2:B:239:GLN:C	2:B:241:VAL:N	2.65	0.52
2:D:73:ARG:NH1	2:D:151:GLN:HE21	2.06	0.52
2:B:21:ASP:OD1	2:B:156:GLN:NE2	2.42	0.52
1:C:8:GLU:OE1	1:C:8:GLU:N	2.41	0.52
1:C:14:ASP:OD1	1:C:14(C):GLU:N	2.42	0.52
2:D:61:GLU:N	2:D:61:GLU:OE1	2.43	0.52
2:D:145:LYS:CG	2:D:149(E):LYS:HD2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:29:TRP:CD2	2:B:121:VAL:HB	2.45	0.52
2:B:61:GLU:CG	2:B:87:LYS:HA	2.38	0.52
2:D:119:HIS:CD2	2:D:120:PRO:HD2	2.44	0.52
2:D:57:HIS:C	2:D:57:HIS:CD2	2.88	0.52
2:B:33:LEU:HD23	2:B:42:CYS:HB2	1.91	0.52
2:D:163:VAL:HB	2:D:182:CYS:SG	2.49	0.52
2:B:101:ARG:O	2:B:103:ILE:N	2.43	0.51
2:D:57:HIS:CD2	2:D:58:CYS:N	2.78	0.51
2:D:204(B):ASN:OD1	2:D:206:ARG:HD2	2.09	0.51
1:A:5:PRO:O	1:A:9:LYS:HB2	2.10	0.51
2:B:75:ARG:NH2	2:B:77:GLU:HG3	2.25	0.51
1:C:4:ARG:CZ	1:C:7:PHE:CD2	2.94	0.51
2:D:61:GLU:OE2	2:D:87:LYS:HA	2.10	0.51
2:D:76:TYR:HE1	2:D:82:ILE:HD11	1.76	0.51
2:D:131:GLN:O	2:D:162:ILE:HD12	2.10	0.51
2:D:149:THR:O	2:D:149(C):VAL:O	2.29	0.51
2:D:66:VAL:HG23	2:D:66:VAL:O	2.10	0.51
2:D:99:LEU:HD12	2:D:99:LEU:H	1.76	0.51
2:B:195:SER:HA	2:B:213:VAL:HB	1.92	0.51
2:D:98:ASN:O	2:D:100:ASP:N	2.43	0.51
2:B:99:LEU:CD2	3:E:46:PRO:HB3	2.40	0.51
2:B:170:ASP:O	2:B:172:THR:N	2.44	0.51
2:B:204(B):ASN:OD1	2:B:206:ARG:CG	2.58	0.51
2:B:239:GLN:O	2:B:242:ILE:N	2.44	0.51
1:C:4:ARG:CZ	1:C:7:PHE:HD2	2.24	0.51
2:D:148:TRP:CE2	2:D:150:GLY:HA2	2.45	0.51
2:B:35:ARG:HB2	2:B:41:LEU:HD21	1.93	0.51
2:B:100:ASP:OD1	2:B:179:ASN:ND2	2.44	0.51
2:D:52:VAL:HG12	2:D:53:LEU:N	2.26	0.51
2:B:232:PHE:O	2:B:233:ARG:C	2.53	0.50
2:D:27:SER:OG	2:D:29:TRP:NE1	2.44	0.50
1:C:14(K):ILE:O	1:C:14(K):ILE:HG22	2.10	0.50
2:D:174:ILE:HD12	2:D:215:TRP:CZ3	2.47	0.50
2:D:29:TRP:HA	2:D:46:LEU:HB3	1.93	0.50
2:D:136:GLY:HA2	2:D:201:MET:HG2	1.93	0.50
2:B:185:LYS:HB3	2:B:186:PRO:CD	2.42	0.50
2:D:164:GLU:O	2:D:167:VAL:N	2.41	0.50
2:B:20:SER:O	2:B:156:GLN:OE1	2.29	0.50
2:B:128:THR:O	2:B:129(B):SER:N	2.45	0.50
2:D:118:ILE:HD12	2:D:118:ILE:N	2.26	0.50
2:D:149(E):LYS:O	2:D:150:GLY:C	2.55	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:197:GLY:O	2:D:212:ILE:HA	2.12	0.50
2:B:57:HIS:C	2:B:57:HIS:CD2	2.89	0.49
2:D:156:GLN:NE2	2:D:156:GLN:HA	2.28	0.49
2:D:197:GLY:O	2:D:212:ILE:HG22	2.11	0.49
1:A:14(G):LEU:HD13	1:A:14(G):LEU:O	2.12	0.49
2:D:146:GLU:O	2:D:147:THR:CB	2.59	0.49
2:D:158:VAL:HG22	2:D:159:ASN:N	2.26	0.49
2:D:136:GLY:HA3	2:D:199:PHE:HE1	1.72	0.49
2:D:60(D):TRP:CH2	3:F:46:PRO:HG2	2.48	0.49
2:D:143:ASN:HD22	2:D:192:GLU:HB2	1.77	0.49
2:B:79:ILE:HG22	2:B:80:GLU:N	2.28	0.49
2:B:164:GLU:OE1	2:B:164:GLU:N	2.40	0.49
1:C:8:GLU:HG3	2:D:205:ASN:ND2	2.27	0.49
2:D:57:HIS:HE2	2:D:195:SER:HB2	1.77	0.49
2:B:57:HIS:HD2	2:B:58:CYS:N	2.11	0.49
2:B:60(D):TRP:O	2:B:60(E):ASP:C	2.55	0.49
2:B:66:VAL:O	2:B:82:ILE:HA	2.12	0.49
2:D:157:VAL:HG22	2:D:158:VAL:N	2.26	0.49
2:D:164:GLU:CD	2:D:164:GLU:H	2.21	0.49
2:D:201:MET:O	2:D:207:TRP:HA	2.13	0.49
2:D:56:ALA:HB1	2:D:94:TYR:CE2	2.47	0.49
2:D:70:LYS:HZ3	2:D:80:GLU:CD	2.21	0.49
2:D:141:TRP:HA	2:D:152:PRO:HG3	1.95	0.49
2:D:243:ASP:C	2:D:245:PHE:H	2.20	0.49
2:B:82:ILE:O	2:B:82:ILE:HG22	2.12	0.48
2:D:49:ASP:CG	2:D:50:ARG:HG2	2.38	0.48
2:D:157:VAL:CG2	2:D:158:VAL:N	2.76	0.48
2:B:50:ARG:HH11	2:B:107:LYS:HE3	1.77	0.48
2:B:168:CYS:O	2:B:170:ASP:N	2.46	0.48
2:D:200:VAL:HG12	2:D:209:GLN:HB2	1.95	0.48
2:B:170:ASP:C	2:B:172:THR:N	2.69	0.48
2:B:230:HIS:O	2:B:231:VAL:C	2.57	0.48
2:D:84:MET:HB3	2:D:109:LYS:NZ	2.28	0.48
2:D:84:MET:O	2:D:108:LEU:HB3	2.14	0.48
2:D:240:LYS:O	2:D:244:GLN:HB3	2.14	0.48
2:B:32:MET:HE2	2:B:40:LEU:HD13	1.96	0.48
2:B:191:CYS:O	2:B:192:GLU:C	2.56	0.48
2:D:45:SER:OG	2:D:53:LEU:HB3	2.13	0.48
2:D:147:THR:HG22	2:D:148:TRP:N	2.27	0.48
2:D:204(B):ASN:ND2	2:D:204(B):ASN:N	2.56	0.48
1:A:14(F):LEU:C	1:A:14(H):GLU:N	2.70	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:192:GLU:N	2:B:192:GLU:OE1	2.46	0.48
2:D:52:VAL:CG1	2:D:53:LEU:N	2.77	0.48
2:D:165:ARG:HB3	2:D:166:PRO:HD3	1.95	0.48
2:D:242:ILE:O	2:D:245:PHE:CE1	2.67	0.48
2:D:202:LYS:O	2:D:204:PRO:HD3	2.13	0.48
2:D:70:LYS:NZ	2:D:70:LYS:HA	2.29	0.48
2:D:94:TYR:HA	2:D:101:ARG:HB2	1.94	0.48
2:D:185:LYS:O	2:D:186(A):ASP:N	2.41	0.48
2:B:127:GLU:HG2	2:D:148:TRP:CH2	2.49	0.47
1:C:13:GLU:HA	1:C:14(C):GLU:OE2	2.13	0.47
1:C:14:ASP:O	1:C:14(B):THR:N	2.47	0.47
2:D:57:HIS:ND1	3:F:46:PRO:CB	2.77	0.47
2:B:86:GLU:OE2	2:B:107:LYS:HE2	2.15	0.47
1:C:14(G):LEU:HD23	1:C:14(J):TYR:CE2	2.49	0.47
2:D:139:THR:HG22	2:D:157:VAL:CB	2.44	0.47
2:D:197:GLY:O	2:D:212:ILE:CG2	2.63	0.47
2:B:130:LEU:O	2:B:130:LEU:HG	2.15	0.47
1:C:1(C):GLU:O	1:C:1(B):ALA:C	2.57	0.47
2:D:211:GLY:HA2	2:D:231:VAL:HG23	1.97	0.47
2:B:38:GLN:O	2:B:38:GLN:OE1	2.32	0.47
2:D:149(B):ASN:C	2:D:149(C):VAL:CG2	2.88	0.47
2:B:41:LEU:CD1	2:B:64:LEU:HD22	2.44	0.47
2:B:203:SER:O	2:B:205:ASN:HA	2.15	0.47
2:D:53:LEU:HD12	2:D:54:THR:N	2.29	0.47
2:D:59:LEU:HD13	2:D:88:ILE:HG21	1.94	0.47
2:D:142:GLY:O	2:D:143:ASN:C	2.58	0.47
2:D:145:LYS:HG2	2:D:149(E):LYS:HD2	1.95	0.47
2:D:70:LYS:HA	2:D:70:LYS:HZ2	1.79	0.47
2:D:179:ASN:O	2:D:230:HIS:HB2	2.15	0.47
2:B:203:SER:HB3	2:B:204(B):ASN:HD21	1.78	0.47
1:C:7:PHE:HD1	1:C:12:LEU:HB3	1.79	0.47
2:D:149(C):VAL:HB	2:D:149(E):LYS:HG3	1.97	0.47
2:B:194:ASP:O	2:B:195:SER:C	2.56	0.47
2:D:195:SER:HA	2:D:213:VAL:CB	2.33	0.47
1:A:1:CYS:O	1:A:3:LEU:N	2.48	0.46
2:B:77(A):ARG:C	2:B:79:ILE:H	2.23	0.46
2:B:80:GLU:C	2:B:81:LYS:HD2	2.40	0.46
2:D:26:MET:HE2	2:D:137:ARG:NH1	2.30	0.46
2:B:114:PHE:CD1	2:B:120:PRO:HD3	2.50	0.46
2:B:131:GLN:HE21	2:B:131:GLN:HB3	1.50	0.46
2:D:239:GLN:O	2:D:242:ILE:N	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:72:SER:HA	2:B:153:SER:O	2.15	0.46
2:B:168:CYS:C	2:B:170:ASP:N	2.73	0.46
2:D:128:THR:HG21	2:D:208:TYR:CD1	2.50	0.46
2:D:169:LYS:HA	2:D:176:ILE:HD13	1.96	0.46
2:B:77:GLU:O	2:B:79:ILE:HB	2.15	0.46
2:D:126:ARG:CZ	2:D:126:ARG:CB	2.94	0.46
2:D:147:THR:O	2:D:148:TRP:CG	2.68	0.46
2:D:32:MET:HE2	2:D:32:MET:HB3	1.85	0.46
2:D:73:ARG:C	2:D:75:ARG:H	2.24	0.46
2:D:219:GLY:HA3	2:D:221(A):ARG:HE	1.81	0.46
2:D:148:TRP:O	2:D:149:THR:OG1	2.28	0.46
2:D:206:ARG:CB	2:D:208:TYR:HE2	2.29	0.46
2:D:60(B):PRO:N	2:D:60(C):PRO:CD	2.79	0.46
2:B:180:MET:HA	2:B:228:TYR:O	2.16	0.46
2:B:73:ARG:NH1	2:B:151:GLN:HG3	2.22	0.45
2:B:100:ASP:OD1	2:B:101:ARG:HG3	2.16	0.45
1:C:3:LEU:CB	1:C:9:LYS:HE3	2.46	0.45
2:D:34:PHE:HE1	2:D:38:GLN:NE2	2.14	0.45
2:D:56:ALA:HB2	2:D:103:ILE:H	1.82	0.45
2:B:86:GLU:O	2:B:86:GLU:HG2	2.15	0.45
1:C:14(G):LEU:C	1:C:14(I):SER:N	2.74	0.45
2:B:32:MET:HE2	2:B:40:LEU:CD1	2.46	0.45
2:B:95:ASN:OD1	2:B:95:ASN:C	2.57	0.45
2:D:50:ARG:HH11	2:D:111:PRO:CD	2.29	0.45
2:D:77:GLU:HB3	2:D:79:ILE:CB	2.47	0.45
2:B:47:ILE:HD11	2:B:51:TRP:CB	2.46	0.45
2:B:195:SER:HB3	3:E:47:ARG:C	2.41	0.45
2:D:77:GLU:CB	2:D:79:ILE:HB	2.46	0.45
2:D:91:HIS:ND1	2:D:93:ARG:N	2.62	0.45
2:B:125:ASP:OD2	2:B:128:THR:HB	2.17	0.45
2:B:204(B):ASN:ND2	2:B:204(B):ASN:H	2.15	0.45
2:D:127:GLU:O	2:D:128:THR:C	2.59	0.45
2:D:57:HIS:CE1	3:F:46:PRO:HB2	2.52	0.45
2:B:23:GLU:O	2:B:26:MET:HB2	2.16	0.45
2:B:57:HIS:CD2	2:B:58:CYS:N	2.84	0.45
2:B:204(B):ASN:OD1	2:B:206:ARG:HB2	2.17	0.45
2:B:233:ARG:C	2:B:235:LYS:H	2.25	0.45
2:B:239:GLN:O	2:B:241:VAL:N	2.50	0.45
1:C:14(G):LEU:C	1:C:14(I):SER:H	2.24	0.45
2:D:148:TRP:NE1	2:D:150:GLY:HA2	2.31	0.45
2:D:147:THR:CG2	2:D:148:TRP:N	2.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:239:GLN:O	2:B:240:LYS:C	2.59	0.45
2:B:20:SER:OG	2:B:21:ASP:N	2.51	0.44
2:D:35:ARG:HD3	2:D:37:PRO:HD2	1.98	0.44
2:B:24:ILE:HG12	2:B:71:HIS:ND1	2.32	0.44
2:B:94:TYR:CZ	2:B:96:TRP:HB3	2.52	0.44
2:B:126:ARG:HB2	2:B:126:ARG:CZ	2.48	0.44
2:B:134:TYR:OH	2:D:149(A):ALA:HB2	2.17	0.44
2:D:96:TRP:CD1	2:D:96:TRP:N	2.85	0.44
2:D:235:LYS:HA	2:D:238:ILE:CG1	2.47	0.44
2:D:108:LEU:N	2:D:108:LEU:HD23	2.32	0.44
2:B:100:ASP:OD1	2:B:101:ARG:N	2.51	0.44
2:B:185:LYS:HB2	2:B:186(B):GLU:HG3	2.00	0.44
2:B:216:GLY:HA3	3:E:47:ARG:NH2	2.33	0.44
2:D:233:ARG:HG2	2:D:233:ARG:HH21	1.82	0.44
2:B:27:SER:O	2:B:30:GLN:HB2	2.18	0.44
2:D:31:VAL:HB	2:D:44:ALA:C	2.43	0.44
2:B:79:ILE:HD13	2:B:79:ILE:HA	1.89	0.44
1:C:4:ARG:NH2	1:C:14(C):GLU:OE2	2.50	0.44
2:D:191:CYS:N	2:D:194:ASP:OD2	2.48	0.44
1:C:14(G):LEU:O	1:C:14(I):SER:N	2.51	0.44
2:D:36:LYS:HD2	2:D:62:ASN:O	2.18	0.44
2:B:35:ARG:CA	2:B:41:LEU:HD11	2.48	0.44
2:B:45:SER:CB	2:B:209:GLN:HE22	2.30	0.44
2:B:93:ARG:HB2	2:B:101:ARG:HD2	2.00	0.44
2:B:129:ALA:HA	2:B:210:MET:HE1	1.99	0.44
2:D:220:CYS:O	2:D:221:ASP:C	2.60	0.44
2:B:123:LEU:HA	2:B:124:PRO:HD3	1.85	0.43
2:B:130:LEU:O	2:B:130:LEU:CG	2.65	0.43
2:D:60(H):PHE:CB	2:D:64:LEU:HD11	2.47	0.43
2:D:96:TRP:CZ2	2:D:97:ARG:HG2	2.53	0.43
2:B:88:ILE:C	2:B:89:TYR:CD1	2.96	0.43
2:D:23:GLU:O	2:D:24:ILE:C	2.61	0.43
2:D:114:PHE:N	2:D:114:PHE:CD1	2.84	0.43
2:B:32:MET:HG2	2:B:141:TRP:CZ3	2.54	0.43
2:B:34:PHE:CB	2:B:65:LEU:HD23	2.49	0.43
2:B:88:ILE:HG22	2:B:89:TYR:N	2.33	0.43
2:B:145:LYS:HG3	2:B:146:GLU:N	2.33	0.43
2:B:163:VAL:HG12	2:B:164:GLU:N	2.32	0.43
2:B:236:LYS:HZ2	2:B:236:LYS:CB	2.29	0.43
1:C:1(A):ASP:OD2	1:C:9:LYS:NZ	2.46	0.43
1:C:7:PHE:CB	1:C:12:LEU:O	2.64	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:204(B):ASN:HD22	2:D:204(B):ASN:N	2.07	0.43
2:D:219:GLY:HA3	2:D:221(A):ARG:NE	2.34	0.43
1:A:6:LEU:HA	1:A:10:LYS:HD3	1.99	0.43
2:D:27:SER:HG	2:D:29:TRP:CD1	2.37	0.43
2:D:45:SER:HB3	2:D:209:GLN:HE22	1.79	0.43
2:B:92:PRO:HD3	2:B:237:TRP:HE1	1.83	0.43
1:C:2:GLY:O	2:D:207:TRP:HD1	2.01	0.43
1:C:14(J):TYR:CD1	2:D:134:TYR:CD2	3.06	0.43
2:D:16:ILE:HD11	2:D:139:THR:CA	2.49	0.43
2:D:108:LEU:O	2:D:109:LYS:C	2.62	0.43
2:B:188:GLY:O	2:B:189:ASP:HB2	2.19	0.43
1:A:14(G):LEU:HD22	1:A:14(G):LEU:HA	1.89	0.43
1:A:14(J):TYR:CD1	2:B:134:TYR:CD2	3.06	0.43
2:B:91:HIS:CG	2:B:92:PRO:HD2	2.54	0.43
2:B:31:VAL:HG12	2:B:32:MET:N	2.34	0.43
2:D:70:LYS:NZ	2:D:77:GLU:OE2	2.51	0.43
2:D:180:MET:HE2	2:D:215:TRP:HE1	1.83	0.43
2:D:238:ILE:HG22	2:D:242:ILE:CD1	2.41	0.43
2:B:56:ALA:N	2:B:102:ASN:OD1	2.52	0.43
2:B:209:GLN:HG2	2:B:231:VAL:HG21	2.00	0.43
2:D:84:MET:CB	2:D:109:LYS:NZ	2.82	0.43
2:D:90:ILE:CG2	2:D:91:HIS:N	2.71	0.43
2:D:91:HIS:HA	2:D:92:PRO:HD3	1.86	0.43
2:D:234:LEU:HD23	2:D:234:LEU:N	2.34	0.43
1:C:14(D):ARG:CD	1:C:14(D):ARG:C	2.91	0.42
2:B:96:TRP:HA	2:B:99:LEU:HD12	2.01	0.42
2:D:35:ARG:O	2:D:35:ARG:HG2	2.19	0.42
2:D:41:LEU:HD11	2:D:60(H):PHE:CZ	2.54	0.42
2:D:70:LYS:HE3	2:D:70:LYS:HB3	1.89	0.42
2:B:94:TYR:HA	2:B:101:ARG:HB2	2.00	0.42
1:C:14(J):TYR:HD1	2:D:134:TYR:CD2	2.38	0.42
2:D:77:GLU:HB3	2:D:79:ILE:CG1	2.49	0.42
2:D:96:TRP:CG	2:D:97:ARG:N	2.87	0.42
1:C:14(F):LEU:HD21	2:D:159:ASN:OD1	2.18	0.42
2:B:29:TRP:HB3	2:B:119:HIS:O	2.20	0.42
2:D:163:VAL:HG12	2:D:164:GLU:N	2.33	0.42
1:C:12:LEU:HD22	1:C:12:LEU:HA	1.80	0.42
1:C:14(G):LEU:HD21	2:D:202:LYS:HG2	2.01	0.42
2:D:129(C):LEU:HD11	2:D:203:SER:HA	2.00	0.42
2:D:199:PHE:CD1	2:D:199:PHE:C	2.97	0.42
2:B:37:PRO:O	2:B:38:GLN:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:100:ASP:OD2	2:B:177:THR:HG21	2.20	0.42
2:D:235:LYS:HG3	2:D:239:GLN:HE21	1.84	0.42
1:C:14(G):LEU:HA	1:C:14(J):TYR:CD2	2.55	0.42
2:D:96:TRP:O	2:D:97(A):GLU:N	2.53	0.42
2:D:195:SER:OG	3:F:47:ARG:O	2.38	0.42
2:D:203:SER:HB3	2:D:204(B):ASN:HD21	1.85	0.42
2:B:47:ILE:CD1	2:B:51:TRP:HB2	2.48	0.41
2:B:122:CYS:HB2	2:B:208:TYR:CD1	2.54	0.41
2:B:75:ARG:HH21	2:B:77:GLU:HA	1.85	0.41
2:B:126:ARG:HA	2:B:232:PHE:CZ	2.55	0.41
2:B:130:LEU:HD11	2:B:181:PHE:CD2	2.55	0.41
2:B:203:SER:C	2:B:204(A):PHE:H	2.27	0.41
2:D:134:TYR:O	2:D:161:PRO:HA	2.21	0.41
2:D:233:ARG:HG2	2:D:233:ARG:NH2	2.35	0.41
2:B:54:THR:HG22	2:B:104:ALA:O	2.20	0.41
2:B:174:ILE:HD11	3:E:44:PRO:HG3	2.02	0.41
2:D:23:GLU:HB2	2:D:26:MET:CG	2.46	0.41
2:D:36(A):SER:HA	2:D:37:PRO:HA	1.74	0.41
2:D:114:PHE:O	2:D:115:SER:CB	2.68	0.41
2:D:212:ILE:O	2:D:214:SER:N	2.53	0.41
2:B:16:ILE:HG21	2:B:158:VAL:CG1	2.50	0.41
2:B:77:GLU:OE2	2:B:77:GLU:N	2.42	0.41
2:B:125:ASP:O	2:B:125:ASP:OD2	2.38	0.41
2:B:239:GLN:HE21	2:B:239:GLN:HB2	1.64	0.41
2:D:65:LEU:HD11	2:D:82:ILE:CG2	2.51	0.41
2:D:151:GLN:HG2	2:D:152:PRO:HD2	2.02	0.41
1:C:14(K):ILE:C	1:C:14(L):ASP:OD1	2.64	0.41
2:B:98:ASN:O	2:B:99:LEU:C	2.63	0.41
2:B:215:TRP:CB	3:E:46:PRO:HA	2.50	0.41
1:C:14(A):LYS:N	2:D:23:GLU:OE2	2.54	0.41
1:C:14(D):ARG:HD3	1:C:14(D):ARG:C	2.45	0.41
2:D:33:LEU:O	2:D:41:LEU:HB2	2.20	0.41
2:D:45:SER:HB2	2:D:209:GLN:HE22	1.81	0.41
2:B:100:ASP:OD1	2:B:101:ARG:CG	2.69	0.41
2:B:112:VAL:HG22	2:B:113:ALA:N	2.36	0.41
1:C:14:ASP:CB	2:D:26:MET:HE3	2.50	0.41
2:D:100:ASP:CG	2:D:177:THR:HG21	2.46	0.41
2:D:238:ILE:O	2:D:242:ILE:HG13	2.21	0.41
2:B:124:PRO:HG3	2:B:210:MET:SD	2.60	0.41
2:D:65:LEU:HD11	2:D:82:ILE:HG21	2.03	0.41
2:D:151:GLN:CG	2:D:152:PRO:HD2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:165:ARG:N	2:B:166:PRO:CD	2.84	0.40
2:B:163:VAL:CG1	2:B:167:VAL:HB	2.50	0.40
2:D:102:ASN:O	2:D:103:ILE:HB	2.21	0.40
2:D:220:CYS:O	2:D:221:ASP:O	2.38	0.40
2:D:187:ARG:HG2	2:D:187:ARG:NH1	2.36	0.40
2:D:208:TYR:O	2:D:210:MET:HG2	2.22	0.40
2:B:95:ASN:HB3	2:B:100:ASP:HB3	2.03	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	27/31 (87%)	18 (67%)	6 (22%)	3 (11%)	0	1
1	C	27/31 (87%)	20 (74%)	3 (11%)	4 (15%)	0	0
2	B	247/259 (95%)	194 (78%)	35 (14%)	18 (7%)	1	2
2	D	255/259 (98%)	194 (76%)	39 (15%)	22 (9%)	0	1
3	E	7/9 (78%)	4 (57%)	0	3 (43%)	0	0
3	F	2/9 (22%)	2 (100%)	0	0	100	100
All	All	565/598 (94%)	432 (76%)	83 (15%)	50 (9%)	0	1

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	GLY
2	B	20	SER
2	B	87	LYS
2	B	102	ASN
2	B	130	LEU

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Mol	Chain	Res	Type
2	B	158	VAL
1	C	14(A)	LYS
2	D	99	LEU
2	D	115	SER
2	D	148	TRP
2	D	149(A)	ALA
2	D	149(B)	ASN
2	D	149(C)	VAL
2	D	186(A)	ASP
2	D	213	VAL
3	E	40	PRO
3	E	42	ILE
1	A	1	CYS
2	B	36	LYS
2	B	60(E)	ASP
2	B	61	GLU
2	B	169	LYS
2	B	195	SER
1	C	1(B)	ALA
2	D	24	ILE
2	D	60(H)	PHE
2	D	97	ARG
2	D	126	ARG
2	D	147	THR
2	D	150	GLY
2	D	195	SER
2	D	221	ASP
2	B	171	SER
2	B	221	ASP
2	D	98	ASN
3	E	41	SER
2	B	54	THR
2	B	75	ARG
1	C	1(A)	ASP
1	C	14(H)	GLU
2	D	114	PHE
2	D	186	PRO
2	B	82	ILE
2	B	129	ALA
2	B	234	LEU
2	B	240	LYS
2	D	127	GLU

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Mol	Chain	Res	Type
1	A	7	PHE
2	D	103	ILE
2	D	231	VAL

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	27/28 (96%)	26 (96%)	1 (4%)	30	65
1	C	27/28 (96%)	26 (96%)	1 (4%)	30	65
2	B	219/225 (97%)	201 (92%)	18 (8%)	10	33
2	D	223/225 (99%)	197 (88%)	26 (12%)	5	18
3	E	8/8 (100%)	6 (75%)	2 (25%)	0	2
3	F	3/8 (38%)	3 (100%)	0	100	100
All	All	507/522 (97%)	459 (90%)	48 (10%)	8	26

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
2	B	41	LEU
2	B	45	SER
2	B	57	HIS
2	B	81	LYS
2	B	99	LEU
2	B	116	ASP
2	B	127	GLU
2	B	128	THR
2	B	131	GLN
2	B	151	GLN
2	B	161	PRO
2	B	173	ARG
2	B	182	CYS
2	B	192	GLU

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Mol	Chain	Res	Type
2	B	199	PHE
2	B	229	THR
2	B	236	LYS
2	B	245	PHE
1	C	12	LEU
2	D	26	MET
2	D	41	LEU
2	D	57	HIS
2	D	63	ASP
2	D	70	LYS
2	D	80	GLU
2	D	81	LYS
2	D	102	ASN
2	D	105	LEU
2	D	108	LEU
2	D	121	VAL
2	D	125	ASP
2	D	147	THR
2	D	151	GLN
2	D	164	GLU
2	D	173	ARG
2	D	176	ILE
2	D	186	PRO
2	D	186(A)	ASP
2	D	186(B)	GLU
2	D	186(D)	LYS
2	D	212	ILE
2	D	224	LYS
2	D	228	TYR
2	D	234	LEU
2	D	236	LYS
3	E	43	LEU
3	E	47	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	91	HIS
2	B	131	GLN
2	B	209	GLN
2	B	239	GLN
2	D	60(G)	ASN

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Mol	Chain	Res	Type
2	D	102	ASN
2	D	143	ASN
2	D	151	GLN
2	D	156	GLN
2	D	179	ASN
2	D	204(B)	ASN
2	D	209	GLN
2	D	230	HIS
2	D	239	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	29/31 (93%)	0.30	1 (3%) 48 39	67, 80, 112, 120	0
1	C	29/31 (93%)	0.43	1 (3%) 48 39	90, 113, 126, 130	0
2	B	251/259 (96%)	0.23	3 (1%) 76 68	43, 78, 102, 116	0
2	D	257/259 (99%)	0.55	5 (1%) 66 57	54, 104, 131, 140	0
3	E	9/9 (100%)	0.85	1 (11%) 10 7	85, 92, 131, 139	0
3	F	4/9 (44%)	1.93	2 (50%) 0 0	126, 137, 140, 141	0
All	All	579/598 (96%)	0.41	13 (2%) 62 52	43, 88, 129, 141	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	106	MET	4.1
3	E	39	THR	3.2
3	F	44	PRO	2.8
2	D	90	ILE	2.8
1	C	1(B)	ALA	2.8
2	B	74	THR	2.6
3	F	45	ALA	2.5
2	D	149	THR	2.5
1	A	1(B)	ALA	2.4
2	B	38	GLN	2.4
2	D	212	ILE	2.3
2	D	105	LEU	2.3
2	B	147	THR	2.1

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.