



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:15 PM UTC

PDB ID : 5KYY / pdb_00005kyx
Title : crystal structure of Sec23 and TANGO1 peptide1 complex
Authors : Ma, W.; Goldberg, J.
Deposited on : 2016-07-22
Resolution : 3.52 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

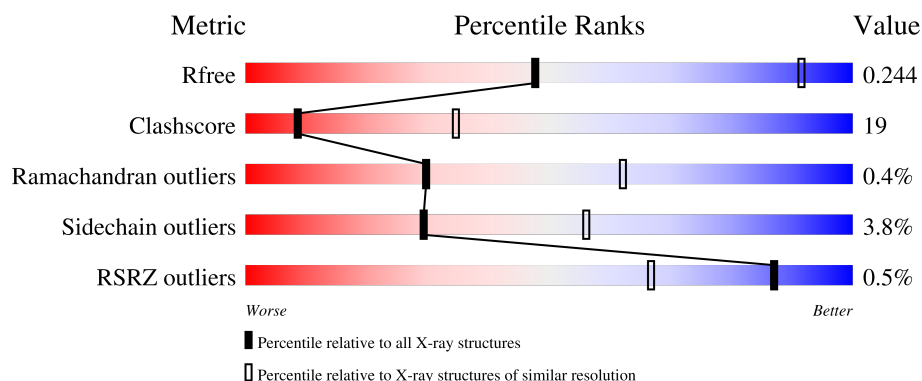
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.52 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	1025 (3.56-3.48)
Clashscore	190562	1079 (3.56-3.48)
Ramachandran outliers	187476	1052 (3.56-3.48)
Sidechain outliers	187428	1053 (3.56-3.48)
RSRZ outliers	180081	1024 (3.56-3.48)

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	765	 67% 23% 6%
2	B	770	 68% 27% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11706 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 Protein transport protein Sec23A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	716	Total	C	N	O	S	0	0	0
			5684	3623	975	1046	40			

- Molecule 2 is a protein called Protein transport protein Sec24D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	767	Total	C	N	O	S	0	0	0
			6020	3831	1017	1118	54			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	ALA	-	expression tag	UNP O94855
B	2	MET	-	expression tag	UNP O94855
B	3	GLY	-	expression tag	UNP O94855

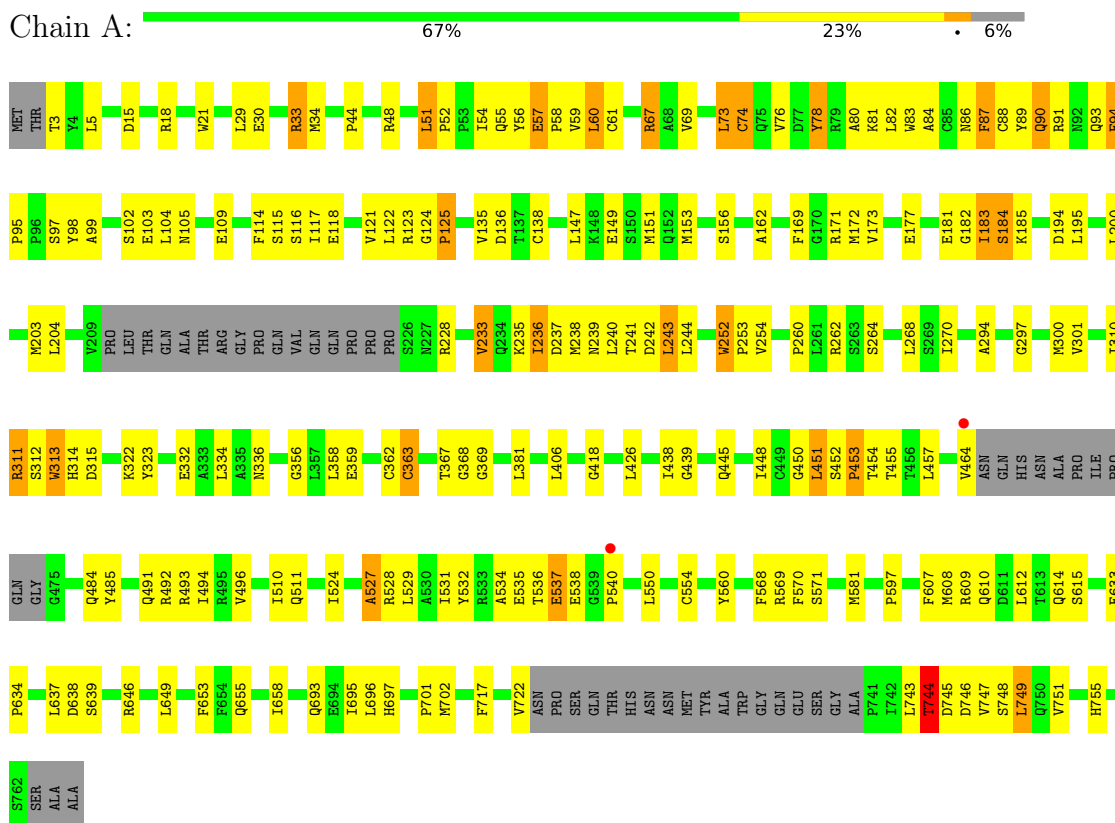
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

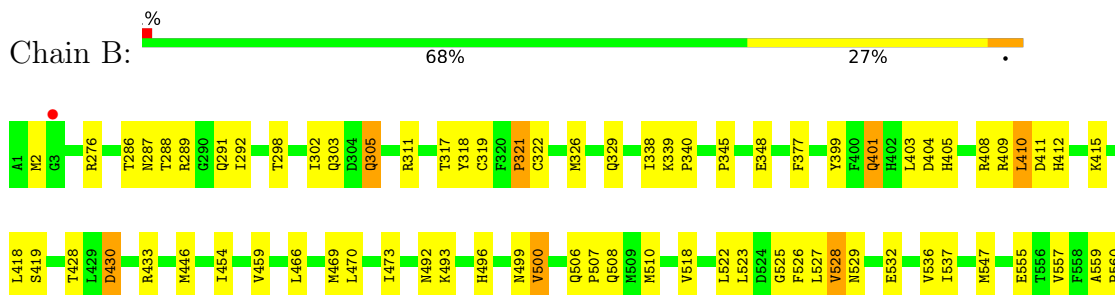
3 Residue-property plots [i](#)

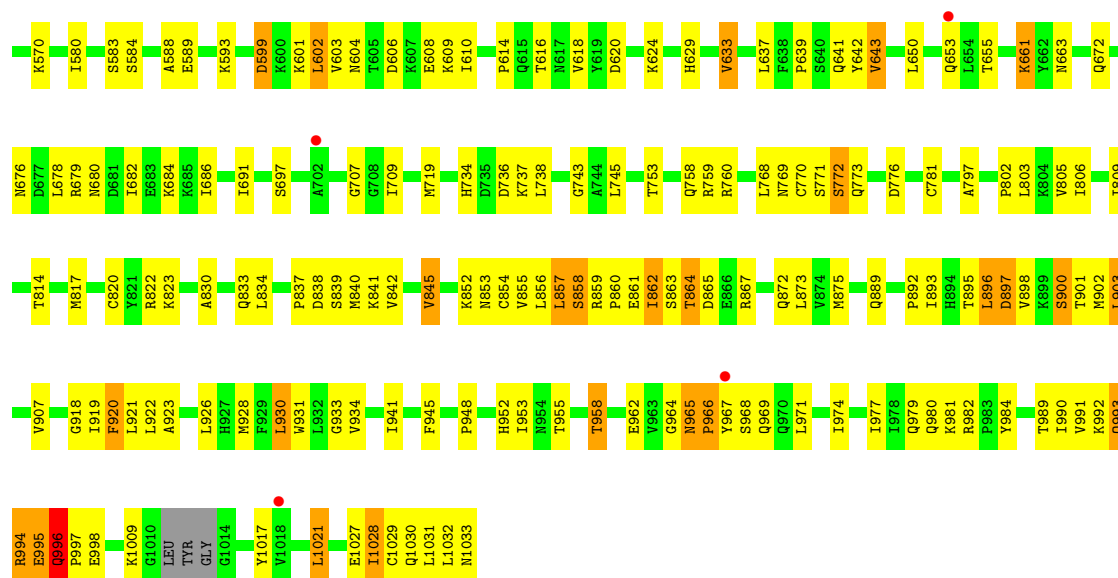
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Protein transport protein Sec23A



- Molecule 2: Protein transport protein Sec24D





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.96Å 141.87Å 152.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.77 – 3.52 48.77 – 3.52	Depositor EDS
% Data completeness (in resolution range)	88.4 (48.77-3.52) 84.1 (48.77-3.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.195 , 0.255 (Not available) , 0.244	Depositor DCC
R_{free} test set	2006 reflections (7.06%)	wwPDB-VP
Wilson B-factor (Å ²)	84.8	Xtriage
Anisotropy	1.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 75.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11706	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.52	0/5817	0.88	21/7878 (0.3%)
2	B	0.52	1/6148 (0.0%)	0.93	31/8330 (0.4%)
All	All	0.52	1/11965 (0.0%)	0.90	52/16208 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	859	ARG	C-N	5.39	1.39	1.33

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	993	GLN	N-CA-C	10.55	122.78	111.28
2	B	858	SER	N-CA-C	9.24	123.00	109.69
2	B	525	GLY	N-CA-C	-8.36	104.05	114.92
1	A	527	ALA	N-CA-C	8.06	120.07	111.28
2	B	857	LEU	N-CA-C	-6.95	98.81	109.65
2	B	772	SER	N-CA-C	-6.90	102.14	111.87
2	B	410	LEU	N-CA-CB	-6.73	106.50	114.17
1	A	252	TRP	CA-C-N	6.32	126.27	119.76
1	A	252	TRP	C-N-CA	6.32	126.27	119.76
1	A	236	ILE	N-CA-C	6.30	118.16	112.17
1	A	744	THR	N-CA-C	6.26	118.05	108.42
2	B	410	LEU	N-CA-C	6.23	123.06	113.28
2	B	996	GLN	CA-C-N	-6.22	112.07	119.84
2	B	996	GLN	C-N-CA	-6.22	112.07	119.84
2	B	599	ASP	N-CA-C	-6.06	101.31	110.28
2	B	862	ILE	N-CA-C	5.96	116.87	108.17
1	A	51	LEU	CA-C-N	-5.94	114.27	120.38
1	A	51	LEU	C-N-CA	-5.94	114.27	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	473	ILE	CA-C-N	5.88	126.22	119.93
2	B	473	ILE	C-N-CA	5.88	126.22	119.93
2	B	859	ARG	CA-C-N	-5.88	113.16	119.98
2	B	859	ARG	C-N-CA	-5.88	113.16	119.98
2	B	994	ARG	N-CA-C	5.87	118.52	109.07
2	B	321	PRO	N-CA-C	-5.75	102.17	111.19
2	B	965	ASN	N-CA-C	5.73	120.02	112.35
1	A	94	PHE	CA-C-N	-5.64	114.58	120.38
1	A	94	PHE	C-N-CA	-5.64	114.58	120.38
2	B	772	SER	CA-C-N	-5.63	115.05	122.99
2	B	772	SER	C-N-CA	-5.63	115.05	122.99
2	B	982	ARG	CA-C-N	5.59	124.95	118.85
2	B	982	ARG	C-N-CA	5.59	124.95	118.85
1	A	57	GLU	CA-C-N	-5.55	115.02	120.52
1	A	57	GLU	C-N-CA	-5.55	115.02	120.52
1	A	184	SER	N-CA-C	5.52	119.21	107.67
2	B	969	GLN	N-CA-C	-5.51	105.28	111.28
2	B	1031	LEU	N-CA-C	-5.48	101.40	109.62
1	A	363	CYS	CA-C-N	-5.46	114.39	120.45
1	A	363	CYS	C-N-CA	-5.46	114.39	120.45
1	A	52	PRO	N-CA-C	-5.45	104.06	110.70
2	B	410	LEU	CB-CA-C	-5.34	102.37	110.06
2	B	857	LEU	CB-CA-C	-5.30	101.46	112.44
1	A	171	ARG	N-CA-C	-5.24	105.56	111.28
1	A	571	SER	N-CA-C	-5.22	103.64	110.53
1	A	81	LYS	CA-C-N	-5.20	114.50	122.62
1	A	81	LYS	C-N-CA	-5.20	114.50	122.62
2	B	864	THR	N-CA-C	-5.18	105.54	111.14
1	A	87	PHE	N-CA-C	5.17	117.00	111.36
2	B	339	LYS	CA-C-N	5.12	124.78	119.56
2	B	339	LYS	C-N-CA	5.12	124.78	119.56
2	B	903	LEU	CA-C-N	-5.07	114.73	119.85
2	B	903	LEU	C-N-CA	-5.07	114.73	119.85
1	A	233	VAL	CB-CA-C	-5.05	105.33	112.14

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	5684	0	5630	225	0
2	B	6020	0	5995	223	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	11706	0	11625	442	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 19.

All (442) 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)
2:B:902:MET:HE3	2:B:903:LEU:H	1.20	1.06
2:B:902:MET:HE3	2:B:903:LEU:N	1.81	0.94
1:A:123:ARG:C	1:A:125:PRO:HD3	1.95	0.91
1:A:528:ARG:HA	1:A:608:MET:HE1	1.53	0.91
1:A:78:TYR:CE2	1:A:102:SER:C	2.51	0.88
1:A:745:ASP:O	1:A:747:VAL:HG23	1.75	0.86
2:B:965:ASN:HB3	2:B:966:PRO:HD3	1.57	0.86
2:B:321:PRO:HA	2:B:326:MET:HE3	1.59	0.85
2:B:736:ASP:OD1	2:B:737:LYS:N	2.10	0.85
1:A:177:GLU:CD	1:A:185:LYS:HE2	2.03	0.83
2:B:493:LYS:HA	2:B:557:VAL:HG13	1.62	0.82
2:B:409:ARG:O	2:B:410:LEU:HB3	1.79	0.82
1:A:747:VAL:HG12	1:A:748:SER:H	1.43	0.81
1:A:749:LEU:HD12	1:A:749:LEU:O	1.79	0.81
1:A:78:TYR:HD2	1:A:102:SER:HA	1.42	0.81
1:A:78:TYR:HE2	1:A:102:SER:C	1.88	0.81
1:A:21:TRP:CE2	1:A:34:MET:CE	2.64	0.81
2:B:930:LEU:HD12	2:B:931:TRP:N	1.96	0.81
2:B:893:ILE:CG2	2:B:971:LEU:HD13	2.10	0.80
1:A:748:SER:OG	1:A:751:VAL:HG23	1.80	0.79
2:B:466:LEU:O	2:B:470:LEU:HG	1.81	0.79
1:A:83:TRP:HD1	1:A:94:PHE:CE2	1.99	0.79
1:A:177:GLU:OE2	1:A:185:LYS:HE2	1.84	0.78
1:A:536:THR:HG22	1:A:537:GLU:H	1.47	0.77
1:A:312:SER:H	1:A:315:ASP:HB2	1.48	0.77
2:B:410:LEU:O	2:B:410:LEU:HD12	1.85	0.77
2:B:955:THR:HG21	2:B:997:PRO:HG3	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:VAL:C	1:A:60:LEU:HD22	2.11	0.76
1:A:149:GLU:O	1:A:153:MET:HG2	1.84	0.76
2:B:902:MET:HE3	2:B:902:MET:CA	2.16	0.76
2:B:852:LYS:NZ	2:B:1009:LYS:O	2.20	0.75
1:A:21:TRP:CD2	1:A:34:MET:HE1	2.22	0.75
1:A:57:GLU:OE1	1:A:58:PRO:HD2	1.87	0.74
1:A:747:VAL:HG12	1:A:748:SER:N	2.03	0.74
1:A:747:VAL:HG11	1:A:751:VAL:HG11	1.68	0.74
2:B:602:LEU:N	2:B:602:LEU:HD12	2.02	0.74
1:A:240:LEU:HD23	1:A:240:LEU:O	1.87	0.74
1:A:311:ARG:HG2	1:A:311:ARG:HH21	1.51	0.74
2:B:893:ILE:HG23	2:B:971:LEU:HD13	1.68	0.74
2:B:923:ALA:HB2	2:B:928:MET:HE2	1.69	0.74
1:A:536:THR:HG22	1:A:537:GLU:N	2.03	0.74
1:A:21:TRP:CZ2	1:A:34:MET:CE	2.71	0.74
1:A:204:LEU:O	1:A:228:ARG:NH2	2.19	0.73
2:B:412:HIS:O	2:B:419:SER:HB3	1.88	0.73
2:B:919:ILE:C	2:B:920:PHE:HD1	1.96	0.73
2:B:633:VAL:H	2:B:655:THR:HG21	1.53	0.73
2:B:326:MET:HE1	2:B:772:SER:HA	1.69	0.73
2:B:599:ASP:OD2	2:B:601:LYS:HE3	1.89	0.73
1:A:312:SER:O	1:A:314:HIS:N	2.22	0.72
1:A:451:LEU:H	1:A:451:LEU:HD12	1.54	0.72
2:B:896:LEU:HD23	2:B:896:LEU:O	1.89	0.72
1:A:183:ILE:HG12	1:A:184:SER:N	2.04	0.72
2:B:919:ILE:C	2:B:920:PHE:CD1	2.68	0.72
2:B:492:ASN:O	2:B:493:LYS:HB3	1.90	0.71
2:B:896:LEU:HD23	2:B:896:LEU:C	2.14	0.71
2:B:492:ASN:ND2	2:B:496:HIS:HE1	1.88	0.71
1:A:78:TYR:CD2	1:A:102:SER:HA	2.25	0.70
2:B:996:GLN:HB3	2:B:997:PRO:HD2	1.72	0.70
1:A:74:CYS:HB3	1:A:84:ALA:O	1.92	0.70
1:A:312:SER:O	1:A:313:TRP:C	2.33	0.70
2:B:922:LEU:C	2:B:922:LEU:HD23	2.17	0.70
2:B:601:LYS:C	2:B:602:LEU:HD12	2.17	0.69
2:B:900:SER:N	2:B:901:THR:HA	2.08	0.69
2:B:930:LEU:HD12	2:B:930:LEU:C	2.17	0.69
2:B:971:LEU:O	2:B:974:ILE:HG12	1.93	0.69
1:A:310:ILE:HG22	1:A:311:ARG:H	1.56	0.69
1:A:747:VAL:HG11	1:A:751:VAL:CG1	2.22	0.69
2:B:902:MET:CE	2:B:903:LEU:H	2.01	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ILE:CG2	1:A:56:TYR:CE2	2.76	0.69
1:A:55:GLN:O	1:A:56:TYR:HB3	1.92	0.68
1:A:747:VAL:HG11	1:A:751:VAL:CB	2.23	0.68
1:A:78:TYR:HD1	1:A:78:TYR:H	1.39	0.68
1:A:311:ARG:NH1	1:A:356:GLY:HA2	2.09	0.68
2:B:858:SER:C	2:B:860:PRO:HD2	2.19	0.68
1:A:311:ARG:HH21	1:A:311:ARG:CG	2.06	0.68
1:A:537:GLU:O	1:A:538:GLU:C	2.33	0.67
2:B:994:ARG:O	2:B:996:GLN:HG2	1.93	0.67
2:B:1029:CYS:O	2:B:1032:LEU:CD2	2.41	0.67
1:A:114:PHE:HB3	1:A:117:ILE:HG21	1.76	0.67
1:A:21:TRP:CH2	1:A:34:MET:HE3	2.29	0.67
2:B:614:PRO:HG3	2:B:650:LEU:HD22	1.77	0.67
1:A:451:LEU:HD12	1:A:451:LEU:N	2.10	0.66
1:A:747:VAL:CG1	1:A:751:VAL:HB	2.25	0.66
1:A:78:TYR:HE2	1:A:102:SER:O	1.78	0.66
2:B:864:THR:O	2:B:867:ARG:N	2.28	0.66
1:A:48:ARG:CG	1:A:51:LEU:HG	2.26	0.66
2:B:900:SER:CB	2:B:901:THR:HA	2.26	0.65
2:B:433:ARG:HH22	2:B:691:ILE:CD1	2.10	0.65
1:A:182:GLY:C	2:B:506:GLN:OE1	2.39	0.65
2:B:321:PRO:CA	2:B:326:MET:HE3	2.26	0.65
2:B:288:THR:OG1	2:B:291:GLN:HB2	1.97	0.65
2:B:322:CYS:H	2:B:326:MET:CE	2.09	0.65
1:A:83:TRP:CD1	1:A:94:PHE:CE2	2.84	0.65
1:A:177:GLU:CD	1:A:185:LYS:CE	2.68	0.65
1:A:310:ILE:HG22	1:A:311:ARG:N	2.12	0.65
2:B:773:GLN:HB2	2:B:776:ASP:OD2	1.97	0.64
1:A:76:VAL:HG12	1:A:78:TYR:CD1	2.32	0.64
1:A:241:THR:HG23	1:A:242:ASP:OD1	1.97	0.64
2:B:902:MET:HE3	2:B:902:MET:HA	1.79	0.64
1:A:48:ARG:HG3	1:A:51:LEU:HG	1.79	0.64
1:A:235:LYS:C	1:A:236:ILE:HD13	2.24	0.63
1:A:312:SER:O	1:A:315:ASP:N	2.30	0.63
2:B:893:ILE:CG2	2:B:971:LEU:CD1	2.76	0.63
2:B:409:ARG:O	2:B:410:LEU:CB	2.44	0.63
2:B:773:GLN:OE1	2:B:773:GLN:HA	1.98	0.63
2:B:287:ASN:OD1	2:B:288:THR:N	2.32	0.63
2:B:900:SER:N	2:B:901:THR:CA	2.61	0.63
1:A:332:GLU:O	1:A:336:ASN:ND2	2.30	0.63
2:B:931:TRP:CZ2	2:B:993:GLN:HG3	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:994:ARG:HG2	2:B:995:GLU:OE2	2.00	0.62
2:B:931:TRP:HZ2	2:B:993:GLN:HG3	1.63	0.62
1:A:78:TYR:HD2	1:A:102:SER:CA	2.12	0.62
2:B:326:MET:HE1	2:B:771:SER:O	1.98	0.62
2:B:345:PRO:HD2	2:B:348:GLU:OE1	2.00	0.62
1:A:21:TRP:CE2	1:A:34:MET:HE1	2.34	0.62
1:A:123:ARG:HB3	1:A:125:PRO:CD	2.30	0.61
2:B:853:ASN:OD1	2:B:855:VAL:HG23	2.01	0.61
1:A:236:ILE:O	1:A:236:ILE:HG22	2.00	0.61
2:B:834:LEU:HD11	2:B:1028:ILE:HD12	1.83	0.61
1:A:747:VAL:HG11	1:A:751:VAL:HB	1.82	0.61
2:B:977:ILE:HA	2:B:980:GLN:HG3	1.83	0.61
2:B:469:MET:HE1	2:B:678:LEU:HG	1.84	0.60
1:A:195:LEU:HD22	1:A:203:MET:HE1	1.83	0.60
1:A:528:ARG:HG2	1:A:608:MET:CE	2.31	0.60
2:B:797:ALA:HB3	2:B:806:ILE:HD13	1.81	0.60
1:A:169:PHE:CE2	1:A:264:SER:HA	2.35	0.60
1:A:531:ILE:O	1:A:534:ALA:N	2.34	0.60
1:A:312:SER:C	1:A:314:HIS:N	2.58	0.59
2:B:601:LYS:O	2:B:602:LEU:HG	2.03	0.59
2:B:609:LYS:O	2:B:610:ILE:C	2.44	0.59
1:A:78:TYR:CD2	1:A:102:SER:CA	2.85	0.59
2:B:845:VAL:HG22	2:B:1017:TYR:CE1	2.37	0.59
1:A:61:CYS:O	1:A:67:ARG:NH1	2.35	0.59
2:B:1029:CYS:O	2:B:1032:LEU:HD21	2.02	0.59
1:A:30:GLU:HA	1:A:33:ARG:HG3	1.84	0.59
1:A:452:SER:O	1:A:454:THR:N	2.36	0.58
1:A:90:GLN:HG3	1:A:91:ARG:N	2.18	0.58
2:B:964:GLY:HA3	2:B:968:SER:HB3	1.85	0.58
2:B:902:MET:HE3	2:B:902:MET:C	2.29	0.58
1:A:123:ARG:CB	1:A:125:PRO:HD3	2.34	0.58
1:A:358:LEU:HD22	1:A:597:PRO:HB3	1.84	0.58
1:A:368:GLY:HA3	1:A:450:GLY:O	2.04	0.58
1:A:177:GLU:OE1	1:A:185:LYS:CE	2.52	0.57
1:A:82:LEU:HD23	1:A:93:GLN:HA	1.85	0.57
2:B:319:CYS:HA	2:B:769:ASN:O	2.04	0.57
1:A:21:TRP:CZ2	1:A:34:MET:HE3	2.40	0.57
1:A:527:ALA:O	1:A:531:ILE:HG13	2.03	0.57
1:A:95:PRO:HG2	1:A:98:TYR:CD1	2.39	0.57
1:A:322:LYS:HE3	1:A:323:TYR:CZ	2.40	0.57
1:A:240:LEU:HD23	1:A:240:LEU:C	2.30	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:893:ILE:HG21	2:B:971:LEU:CD1	2.35	0.57
2:B:492:ASN:ND2	2:B:496:HIS:CE1	2.72	0.56
2:B:948:PRO:HG2	2:B:952:HIS:HD1	1.70	0.56
1:A:78:TYR:CD1	1:A:78:TYR:N	2.72	0.56
1:A:240:LEU:CD2	1:A:244:LEU:HG	2.36	0.56
2:B:322:CYS:SG	2:B:771:SER:C	2.88	0.56
2:B:920:PHE:CD1	2:B:920:PHE:N	2.72	0.56
1:A:484:GLN:HG2	1:A:494:ILE:HG12	1.87	0.56
1:A:607:PHE:HD2	1:A:608:MET:HE2	1.70	0.56
2:B:922:LEU:HD23	2:B:923:ALA:N	2.20	0.56
2:B:433:ARG:HH22	2:B:691:ILE:HD12	1.69	0.56
2:B:603:VAL:HG12	2:B:604:ASN:OD1	2.06	0.56
2:B:897:ASP:HB3	2:B:900:SER:OG	2.05	0.56
2:B:941:ILE:HG23	2:B:953:ILE:HD11	1.87	0.56
2:B:641:GLN:HG2	2:B:642:TYR:H	1.71	0.55
2:B:1032:LEU:O	2:B:1033:ASN:HB2	2.06	0.55
1:A:241:THR:HG23	1:A:242:ASP:N	2.21	0.55
2:B:893:ILE:HG23	2:B:971:LEU:CD1	2.36	0.55
1:A:749:LEU:HD12	1:A:749:LEU:C	2.27	0.55
2:B:945:PHE:HE2	2:B:990:ILE:HD13	1.71	0.55
2:B:510:MET:SD	2:B:522:LEU:HG	2.47	0.55
2:B:820:CYS:HA	2:B:823:LYS:HE2	1.87	0.55
2:B:979:GLN:NE2	2:B:984:TYR:O	2.40	0.55
1:A:268:LEU:HG	1:A:334:LEU:HD13	1.89	0.54
1:A:560:TYR:HB3	1:A:568:PHE:HA	1.90	0.54
1:A:76:VAL:CG1	1:A:78:TYR:CE1	2.91	0.54
1:A:181:GLU:OE2	1:A:239:ASN:OD1	2.24	0.54
2:B:822:ARG:NH2	2:B:830:ALA:O	2.40	0.54
1:A:123:ARG:HB3	1:A:125:PRO:HD3	1.90	0.54
2:B:977:ILE:HA	2:B:980:GLN:CD	2.33	0.54
1:A:172:MET:HE2	1:A:252:TRP:HE1	1.73	0.54
1:A:88:CYS:O	1:A:89:TYR:HB2	2.07	0.54
1:A:78:TYR:HD1	1:A:78:TYR:N	2.04	0.54
1:A:183:ILE:HG13	2:B:507:PRO:O	2.07	0.54
2:B:966:PRO:HG2	2:B:967:TYR:H	1.72	0.54
2:B:852:LYS:O	2:B:857:LEU:HD22	2.07	0.53
2:B:902:MET:CA	2:B:902:MET:CE	2.87	0.53
1:A:568:PHE:CD1	1:A:568:PHE:C	2.85	0.53
2:B:2:MET:HG2	2:B:1017:TYR:OH	2.08	0.53
1:A:123:ARG:HB3	1:A:125:PRO:HD2	1.90	0.53
1:A:147:LEU:HG	1:A:151:MET:HE2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:554:CYS:HG	1:A:570:PHE:HE2	1.54	0.53
1:A:30:GLU:O	1:A:34:MET:HG3	2.08	0.53
2:B:492:ASN:HD21	2:B:496:HIS:CE1	2.27	0.53
2:B:637:LEU:HD21	2:B:643:VAL:HG11	1.88	0.53
1:A:426:LEU:HD12	1:A:445:GLN:HB3	1.90	0.53
2:B:902:MET:CE	2:B:902:MET:HA	2.37	0.53
1:A:311:ARG:NE	1:A:359:GLU:OE2	2.41	0.53
1:A:21:TRP:CE2	1:A:34:MET:HE2	2.40	0.53
1:A:82:LEU:CD2	1:A:93:GLN:HA	2.38	0.53
1:A:745:ASP:CG	1:A:746:ASP:H	2.17	0.53
2:B:606:ASP:C	2:B:608:GLU:H	2.17	0.53
2:B:864:THR:O	2:B:865:ASP:C	2.50	0.53
1:A:452:SER:C	1:A:454:THR:H	2.17	0.53
1:A:124:GLY:N	1:A:125:PRO:HD3	2.24	0.53
1:A:536:THR:CG2	1:A:537:GLU:N	2.72	0.53
2:B:672:GLN:NE2	2:B:676:ASN:OD1	2.42	0.53
1:A:200:LEU:HD21	1:A:270:ILE:HG23	1.92	0.52
1:A:78:TYR:CD2	1:A:102:SER:C	2.87	0.52
2:B:415:LYS:HB2	2:B:418:LEU:HD12	1.92	0.52
2:B:697:SER:HB3	2:B:745:LEU:HB2	1.92	0.52
2:B:994:ARG:O	2:B:995:GLU:C	2.53	0.52
2:B:965:ASN:HB3	2:B:966:PRO:CD	2.35	0.52
2:B:996:GLN:HB3	2:B:997:PRO:CD	2.39	0.52
1:A:177:GLU:OE1	1:A:185:LYS:HE3	2.08	0.52
2:B:321:PRO:HG2	2:B:734:HIS:HE1	1.75	0.52
2:B:322:CYS:N	2:B:326:MET:CE	2.73	0.52
1:A:238:MET:O	1:A:241:THR:HG22	2.09	0.52
2:B:302:ILE:HB	2:B:348:GLU:OE2	2.10	0.52
2:B:411:ASP:N	2:B:411:ASP:OD1	2.42	0.52
1:A:136:ASP:HA	1:A:169:PHE:CE1	2.45	0.52
1:A:60:LEU:CD2	1:A:60:LEU:N	2.72	0.51
1:A:311:ARG:HE	1:A:359:GLU:CD	2.19	0.51
2:B:318:TYR:HA	2:B:768:LEU:HD22	1.91	0.51
2:B:934:VAL:HG22	2:B:993:GLN:OE1	2.10	0.51
1:A:15:ASP:OD1	1:A:115:SER:OG	2.17	0.51
1:A:264:SER:HB2	1:A:294:ALA:HB2	1.92	0.51
2:B:322:CYS:H	2:B:326:MET:HE1	1.76	0.51
2:B:303:GLN:OE1	2:B:305:GLN:NE2	2.44	0.50
2:B:321:PRO:HA	2:B:326:MET:CE	2.38	0.50
2:B:433:ARG:HH22	2:B:691:ILE:HD11	1.77	0.50
1:A:491:GLN:OE1	1:A:493:ARG:NH2	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:TRP:CZ2	1:A:34:MET:HE2	2.45	0.50
2:B:920:PHE:HD1	2:B:920:PHE:N	2.06	0.50
2:B:922:LEU:C	2:B:922:LEU:CD2	2.85	0.50
1:A:48:ARG:HG2	1:A:51:LEU:HG	1.92	0.50
1:A:452:SER:C	1:A:454:THR:N	2.68	0.50
2:B:934:VAL:HG13	2:B:993:GLN:OE1	2.12	0.50
1:A:3:THR:HG22	1:A:5:LEU:H	1.76	0.50
1:A:252:TRP:CZ3	2:B:523:LEU:HB2	2.47	0.50
2:B:862:ILE:HG22	2:B:863:SER:O	2.11	0.50
1:A:369:GLY:O	1:A:609:ARG:NH2	2.45	0.49
2:B:971:LEU:HD12	2:B:974:ILE:HD11	1.92	0.49
2:B:639:PRO:HB3	2:B:643:VAL:HG21	1.93	0.49
2:B:873:LEU:HD13	2:B:926:LEU:HD11	1.93	0.49
2:B:322:CYS:SG	2:B:771:SER:O	2.70	0.49
2:B:601:LYS:C	2:B:602:LEU:CD1	2.85	0.49
1:A:116:SER:C	1:A:117:ILE:HG23	2.36	0.49
1:A:535:GLU:OE1	1:A:535:GLU:HA	2.13	0.49
2:B:965:ASN:CB	2:B:966:PRO:HD3	2.38	0.49
1:A:78:TYR:CE2	1:A:103:GLU:N	2.79	0.49
1:A:240:LEU:C	1:A:240:LEU:CD2	2.85	0.49
1:A:243:LEU:C	1:A:243:LEU:CD2	2.85	0.49
1:A:746:ASP:C	1:A:747:VAL:CG2	2.85	0.49
2:B:601:LYS:C	2:B:602:LEU:CG	2.85	0.49
2:B:873:LEU:HD13	2:B:926:LEU:CD1	2.43	0.49
1:A:653:PHE:HE2	1:A:701:PRO:HG2	1.78	0.49
2:B:291:GLN:HG2	2:B:292:ILE:N	2.28	0.49
2:B:771:SER:OG	2:B:776:ASP:OD2	2.21	0.49
2:B:918:GLY:HA3	2:B:920:PHE:HE1	1.78	0.49
1:A:116:SER:C	1:A:117:ILE:CG2	2.86	0.49
2:B:977:ILE:HA	2:B:980:GLN:CG	2.43	0.48
2:B:637:LEU:HG	2:B:639:PRO:HD3	1.95	0.48
2:B:996:GLN:C	2:B:998:GLU:H	2.20	0.48
1:A:363:CYS:O	1:A:367:THR:OG1	2.27	0.48
1:A:745:ASP:CG	1:A:746:ASP:N	2.72	0.48
2:B:680:ASN:O	2:B:684:LYS:HG3	2.14	0.48
1:A:252:TRP:HA	1:A:253:PRO:HD3	1.70	0.48
1:A:311:ARG:CG	1:A:311:ARG:NH2	2.71	0.48
1:A:568:PHE:CD1	1:A:569:ARG:N	2.82	0.48
2:B:889:GLN:O	2:B:922:LEU:HA	2.13	0.48
2:B:948:PRO:HG2	2:B:952:HIS:ND1	2.27	0.48
1:A:183:ILE:HG12	1:A:184:SER:H	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:537:GLU:H	1:A:537:GLU:HG3	1.47	0.48
1:A:69:VAL:HG11	1:A:496:VAL:HG21	1.96	0.48
1:A:746:ASP:O	1:A:747:VAL:HG22	2.14	0.48
1:A:124:GLY:N	1:A:125:PRO:CD	2.77	0.48
1:A:182:GLY:O	2:B:506:GLN:OE1	2.32	0.48
1:A:536:THR:CG2	1:A:537:GLU:H	2.19	0.48
1:A:76:VAL:HG12	1:A:78:TYR:CE1	2.49	0.47
2:B:286:THR:O	2:B:305:GLN:O	2.31	0.47
1:A:18:ARG:CZ	1:A:612:LEU:HD22	2.44	0.47
1:A:114:PHE:HB3	1:A:117:ILE:CG2	2.43	0.47
2:B:1027:GLU:HA	2:B:1030:GLN:HB2	1.97	0.47
2:B:854:CYS:O	2:B:867:ARG:HD2	2.13	0.47
2:B:854:CYS:HA	2:B:857:LEU:HB2	1.97	0.47
2:B:428:THR:OG1	2:B:430:ASP:OD1	2.22	0.47
1:A:528:ARG:NH2	1:A:608:MET:C	2.72	0.47
1:A:747:VAL:CG1	1:A:751:VAL:CB	2.86	0.47
2:B:583:SER:OG	2:B:584:SER:N	2.47	0.47
1:A:54:ILE:HG23	1:A:56:TYR:CE2	2.50	0.47
1:A:59:VAL:C	1:A:60:LEU:CD2	2.85	0.47
2:B:377:PHE:HB3	2:B:403:LEU:HD21	1.97	0.47
2:B:399:TYR:O	2:B:409:ARG:NH2	2.48	0.47
2:B:446:MET:HE3	2:B:580:ILE:HG12	1.97	0.47
2:B:499:ASN:HB3	2:B:508:GLN:HB2	1.96	0.47
1:A:237:ASP:O	1:A:241:THR:HG22	2.14	0.47
2:B:289:ARG:O	2:B:769:ASN:ND2	2.48	0.47
2:B:410:LEU:O	2:B:410:LEU:CD1	2.59	0.47
2:B:814:THR:HA	2:B:817:MET:HE2	1.96	0.46
1:A:511:GLN:CD	1:A:511:GLN:H	2.24	0.46
2:B:345:PRO:HG2	2:B:348:GLU:OE1	2.15	0.46
2:B:803:LEU:HD13	2:B:856:LEU:HA	1.97	0.46
2:B:838:ASP:HA	2:B:841:LYS:HG3	1.96	0.46
2:B:931:TRP:CH2	2:B:933:GLY:HA2	2.50	0.46
1:A:747:VAL:CG1	1:A:748:SER:H	2.20	0.46
2:B:861:GLU:O	2:B:861:GLU:HG2	2.15	0.46
1:A:97:SER:C	1:A:99:ALA:H	2.22	0.46
1:A:301:VAL:HG21	1:A:356:GLY:HA3	1.97	0.46
1:A:655:GLN:HE22	1:A:717:PHE:HB3	1.81	0.46
2:B:454:ILE:HD13	2:B:459:VAL:HG21	1.98	0.46
2:B:892:PRO:HB2	2:B:895:THR:HG22	1.98	0.46
1:A:54:ILE:HG21	1:A:56:TYR:CZ	2.50	0.46
2:B:900:SER:HB2	2:B:901:THR:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:709:ILE:HG22	2:B:719:MET:HG2	1.98	0.46
2:B:492:ASN:O	2:B:493:LYS:CB	2.57	0.45
2:B:606:ASP:OD1	2:B:802:PRO:HG3	2.16	0.45
2:B:857:LEU:HD12	2:B:857:LEU:HA	1.77	0.45
1:A:95:PRO:CG	1:A:98:TYR:CD1	2.99	0.45
1:A:310:ILE:HG22	1:A:311:ARG:HG2	1.99	0.45
2:B:834:LEU:CD1	2:B:1028:ILE:CD1	2.95	0.45
2:B:919:ILE:HG23	2:B:930:LEU:HD11	1.98	0.45
1:A:452:SER:O	1:A:453:PRO:C	2.59	0.45
2:B:834:LEU:CD1	2:B:1028:ILE:HD12	2.47	0.45
1:A:54:ILE:HG22	1:A:56:TYR:CD2	2.52	0.45
1:A:121:VAL:HB	1:A:492:ARG:HB3	1.99	0.45
1:A:439:GLY:HA2	1:A:532:TYR:CZ	2.52	0.45
1:A:531:ILE:O	1:A:532:TYR:C	2.57	0.45
2:B:492:ASN:HD21	2:B:496:HIS:HE1	1.60	0.45
1:A:649:LEU:HD23	1:A:658:ILE:HG12	1.99	0.45
1:A:73:LEU:HB2	1:A:86:ASN:HD22	1.81	0.45
1:A:86:ASN:OD1	1:A:87:PHE:CD2	2.70	0.45
1:A:138:CYS:HB2	1:A:262:ARG:NH1	2.32	0.45
2:B:962:GLU:C	2:B:962:GLU:CD	2.85	0.45
1:A:297:GLY:H	1:A:300:MET:CE	2.29	0.44
2:B:559:ALA:HB3	2:B:560:PRO:HD3	2.00	0.44
1:A:136:ASP:CA	1:A:169:PHE:CE1	3.00	0.44
1:A:173:VAL:HG11	1:A:270:ILE:HD12	1.99	0.44
2:B:822:ARG:NH1	2:B:833:GLN:O	2.49	0.44
2:B:921:LEU:HD23	2:B:921:LEU:HA	1.80	0.44
1:A:236:ILE:HD13	1:A:236:ILE:N	2.29	0.44
2:B:837:PRO:HG2	2:B:840:MET:HB2	1.99	0.44
1:A:183:ILE:CG1	1:A:184:SER:N	2.78	0.44
1:A:747:VAL:HG12	1:A:751:VAL:HB	2.00	0.44
2:B:547:MET:HE3	2:B:547:MET:HB2	1.91	0.44
2:B:653:GLN:CD	2:B:872:GLN:HG3	2.43	0.44
2:B:958:THR:HG21	2:B:989:THR:HA	1.99	0.44
1:A:743:LEU:O	1:A:744:THR:HB	2.17	0.44
2:B:326:MET:O	2:B:329:GLN:HB2	2.18	0.44
1:A:297:GLY:H	1:A:300:MET:HE3	1.83	0.44
1:A:638:ASP:OD1	1:A:639:SER:N	2.51	0.44
2:B:805:VAL:O	2:B:809:ILE:HG13	2.18	0.43
2:B:555:GLU:HB3	2:B:588:ALA:HB2	1.99	0.43
2:B:992:LYS:HB2	2:B:996:GLN:HG3	2.00	0.43
1:A:21:TRP:CD2	1:A:34:MET:CE	2.92	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:TYR:OH	1:A:109:GLU:OE2	2.20	0.43
1:A:181:GLU:CG	1:A:239:ASN:OD1	2.66	0.43
2:B:743:GLY:HA2	2:B:770:CYS:HB3	2.01	0.43
1:A:48:ARG:CZ	1:A:51:LEU:HD21	2.48	0.43
2:B:601:LYS:C	2:B:602:LEU:HG	2.43	0.43
2:B:287:ASN:OD1	2:B:288:THR:HG23	2.18	0.43
2:B:559:ALA:N	2:B:560:PRO:CD	2.81	0.43
2:B:758:GLN:OE1	2:B:760:ARG:NH2	2.51	0.43
2:B:276:ARG:HA	2:B:298:THR:HG23	2.01	0.43
2:B:992:LYS:HD2	2:B:996:GLN:HE21	1.83	0.43
1:A:310:ILE:CG2	1:A:311:ARG:N	2.82	0.43
2:B:661:LYS:HE2	2:B:663:ASN:OD1	2.19	0.43
2:B:992:LYS:CB	2:B:996:GLN:HG3	2.49	0.43
1:A:29:LEU:O	1:A:33:ARG:HG2	2.19	0.42
1:A:610:GLN:HB3	1:A:614:GLN:HB2	2.00	0.42
2:B:317:THR:HG22	2:B:319:CYS:H	1.84	0.42
2:B:753:THR:OG1	2:B:759:ARG:NH1	2.49	0.42
1:A:243:LEU:O	1:A:244:LEU:C	2.59	0.42
2:B:624:LYS:HB3	2:B:624:LYS:HE2	1.82	0.42
1:A:153:MET:O	1:A:156:SER:OG	2.37	0.42
1:A:448:ILE:HD11	1:A:457:LEU:HD11	2.00	0.42
1:A:746:ASP:C	1:A:747:VAL:HG23	2.43	0.42
2:B:857:LEU:O	2:B:867:ARG:NH1	2.52	0.42
2:B:991:VAL:HG12	2:B:992:LYS:N	2.33	0.42
1:A:183:ILE:HA	2:B:506:GLN:CD	2.45	0.42
2:B:529:ASN:OD1	2:B:532:GLU:HG2	2.18	0.42
2:B:781:CYS:O	2:B:839:SER:HB2	2.20	0.42
1:A:135:VAL:O	1:A:169:PHE:HD1	2.02	0.42
1:A:359:GLU:OE1	1:A:359:GLU:N	2.52	0.42
1:A:381:LEU:HD12	1:A:702:MET:HE2	2.01	0.42
2:B:526:PHE:O	2:B:527:LEU:HB2	2.20	0.42
1:A:633:GLU:HA	1:A:634:PRO:HD3	1.94	0.42
1:A:743:LEU:O	1:A:755:HIS:ND1	2.52	0.42
2:B:404:ASP:HB3	2:B:405:HIS:H	1.72	0.42
1:A:44:PRO:HD2	1:A:455:THR:O	2.20	0.42
2:B:408:ARG:HD2	2:B:408:ARG:HA	1.83	0.42
2:B:967:TYR:O	2:B:971:LEU:N	2.52	0.42
1:A:104:LEU:HD23	1:A:104:LEU:HA	1.89	0.42
1:A:528:ARG:HG2	1:A:608:MET:HE1	2.02	0.42
2:B:897:ASP:OD1	2:B:898:VAL:N	2.53	0.42
2:B:322:CYS:HA	2:B:738:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:524:ILE:HD12	1:A:615:SER:HB3	2.00	0.41
1:A:58:PRO:HG2	1:A:60:LEU:HD21	2.01	0.41
1:A:76:VAL:HG11	1:A:78:TYR:CE1	2.55	0.41
1:A:73:LEU:HD23	1:A:73:LEU:N	2.35	0.41
1:A:550:LEU:HD13	1:A:581:MET:HG3	2.02	0.41
1:A:607:PHE:CD2	1:A:608:MET:HE2	2.52	0.41
2:B:616:THR:OG1	2:B:618:VAL:HG22	2.20	0.41
1:A:162:ALA:O	1:A:233:VAL:HG23	2.20	0.41
1:A:693:GLN:O	1:A:697:HIS:ND1	2.53	0.41
2:B:322:CYS:N	2:B:326:MET:HE2	2.36	0.41
2:B:500:VAL:O	2:B:528:VAL:HG21	2.20	0.41
1:A:528:ARG:CA	1:A:608:MET:HE1	2.38	0.41
1:A:554:CYS:SG	1:A:570:PHE:HE2	2.42	0.41
2:B:897:ASP:OD1	2:B:900:SER:OG	2.34	0.41
1:A:169:PHE:HE2	1:A:264:SER:HA	1.82	0.41
1:A:254:VAL:HG21	1:A:260:PRO:HG3	2.02	0.41
1:A:312:SER:C	1:A:314:HIS:H	2.28	0.41
2:B:928:MET:HB2	2:B:928:MET:HE3	1.66	0.41
1:A:381:LEU:HA	1:A:702:MET:HE2	2.02	0.41
1:A:118:GLU:OE2	1:A:485:TYR:OH	2.21	0.41
1:A:406:LEU:O	1:A:445:GLN:HA	2.21	0.41
2:B:500:VAL:HG11	2:B:537:ILE:HG12	2.03	0.41
2:B:834:LEU:HD23	2:B:1021:LEU:HD22	2.02	0.41
1:A:80:ALA:O	1:A:82:LEU:HG	2.21	0.41
1:A:418:GLY:HA3	1:A:438:ILE:O	2.21	0.41
2:B:311:ARG:HH12	2:B:981:LYS:NZ	2.19	0.41
2:B:338:ILE:O	2:B:340:PRO:HD3	2.21	0.41
2:B:589:GLU:HA	2:B:593:LYS:HD2	2.03	0.41
2:B:707:GLY:HA2	2:B:875:MET:O	2.21	0.41
1:A:194:ASP:OD1	1:A:195:LEU:N	2.54	0.40
1:A:238:MET:O	1:A:238:MET:SD	2.79	0.40
2:B:401:GLN:HB2	2:B:409:ARG:NH2	2.36	0.40
2:B:469:MET:HE2	2:B:682:ILE:HD12	2.02	0.40
2:B:994:ARG:HG2	2:B:995:GLU:CD	2.45	0.40
1:A:745:ASP:OD2	1:A:746:ASP:N	2.54	0.40
2:B:797:ALA:CB	2:B:806:ILE:HD13	2.51	0.40
1:A:54:ILE:CG2	1:A:56:TYR:CD2	3.04	0.40
1:A:238:MET:SD	1:A:242:ASP:OD1	2.79	0.40
1:A:747:VAL:CG1	1:A:751:VAL:HG21	2.52	0.40
2:B:469:MET:HE3	2:B:679:ARG:HA	2.03	0.40
2:B:570:LYS:HE2	2:B:629:HIS:NE2	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:958:THR:HG23	2:B:990:ILE:HG13	2.03	0.40
1:A:182:GLY:C	2:B:506:GLN:CD	2.90	0.40
2:B:321:PRO:CA	2:B:326:MET:CE	2.96	0.40
2:B:493:LYS:HG2	2:B:493:LYS:O	2.21	0.40
2:B:1032:LEU:O	2:B:1033:ASN:CB	2.69	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	708/765 (92%)	666 (94%)	38 (5%)	4 (1%)	21	55
2	B	763/770 (99%)	703 (92%)	58 (8%)	2 (0%)	36	66
All	All	1471/1535 (96%)	1369 (93%)	96 (6%)	6 (0%)	30	62

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	966	PRO
1	A	313	TRP
2	B	996	GLN
1	A	453	PRO
1	A	540	PRO
1	A	125	PRO

5.3.2 Protein sidechains [i](#)

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	626/666 (94%)	601 (96%)	25 (4%)	28	54
2	B	676/682 (99%)	651 (96%)	25 (4%)	30	57
All	All	1302/1348 (97%)	1252 (96%)	50 (4%)	29	56

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	A	60	LEU
1	A	67	ARG
1	A	73	LEU
1	A	74	CYS
1	A	78	TYR
1	A	90	GLN
1	A	105	ASN
1	A	122	LEU
1	A	183	ILE
1	A	243	LEU
1	A	311	ARG
1	A	362	CYS
1	A	451	LEU
1	A	464	VAL
1	A	510	ILE
1	A	529	LEU
1	A	537	GLU
1	A	637	LEU
1	A	646	ARG
1	A	695	ILE
1	A	696	LEU
1	A	722	VAL
1	A	744	THR
1	A	749	LEU
2	B	305	GLN
2	B	401	GLN
2	B	430	ASP
2	B	500	VAL
2	B	518	VAL
2	B	528	VAL
2	B	536	VAL
2	B	602	LEU

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Mol	Chain	Res	Type
2	B	620	ASP
2	B	633	VAL
2	B	643	VAL
2	B	661	LYS
2	B	686	ILE
2	B	842	VAL
2	B	845	VAL
2	B	896	LEU
2	B	897	ASP
2	B	900	SER
2	B	907	VAL
2	B	920	PHE
2	B	930	LEU
2	B	958	THR
2	B	995	GLU
2	B	1021	LEU
2	B	1028	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	A	90	GLN
1	A	126	GLN
1	A	152	GLN
1	A	227	ASN
1	A	314	HIS
1	A	427	ASN
1	A	436	ASN
1	A	591	GLN
1	A	620	GLN
1	A	710	HIS
2	B	412	HIS
2	B	492	ASN
2	B	496	HIS
2	B	734	HIS
2	B	788	ASN
2	B	833	GLN
2	B	996	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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	716/765 (93%)	-0.19	2 (0%)	90 75	53, 98, 161, 245	0
2	B	767/770 (99%)	-0.15	5 (0%)	84 61	50, 103, 180, 251	0
All	All	1483/1535 (96%)	-0.17	7 (0%)	87 67	50, 101, 171, 251	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	702	ALA	2.8
2	B	967	TYR	2.6
2	B	3	GLY	2.2
2	B	1018	VAL	2.2
1	A	540	PRO	2.1
2	B	653	GLN	2.1
1	A	464	VAL	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	B	1101	1/1	0.95	0.09	90,90,90,90	0
3	ZN	A	801	1/1	0.99	0.06	79,79,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.