Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 03:06 PM GMT

PDB ID	:	4BJL
Title	:	LOCW, A LAMBDA 1 TYPE LIGHT-CHAIN DIMER (BENCE-JONES PROTEIN) CRYSTALLIZED IN DISTILLED WATER
Authors	:	Schiffer, M.; Huang, D.B.
Deposited on	:	1995-05-26
Resolution	:	2.40 Å(reported)

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

The following versions of software and data (see references (1)) were used in the production of this report:

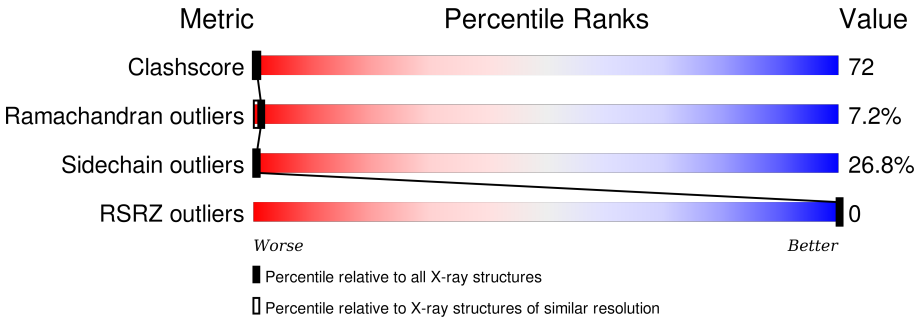
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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	216	22% 44% 30% 5%
1	B	216	19% 50% 25% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3460 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 LOC - LAMBDA 1 TYPE LIGHT-CHAIN DIMER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1598	992	266	335	5			
1	B	216	Total	C	N	O	S	0	0	0
			1598	992	266	335	5			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	THR	ILE	CONFLICT	PIR S25754
A	31	GLU	GLY	CONFLICT	PIR S25754
A	33	SER	THR	CONFLICT	PIR S25754
A	35	THR	ASN	CONFLICT	PIR S25754
A	39	HIS	GLN	CONFLICT	PIR S25754
A	41	SER	PRO	CONFLICT	PIR S25754
A	43	THR	ARG	CONFLICT	PIR S25754
A	50	TYR	HIS	CONFLICT	PIR S25754
A	51	GLU	SER	CONFLICT	PIR S25754
A	52	ASP	ASN	CONFLICT	PIR S25754
A	54	SER	GLN	CONFLICT	PIR S25754
A	56	ALA	PRO	CONFLICT	PIR S25754
A	60	SER	PRO	CONFLICT	PIR S25754
A	65	ALA	GLY	CONFLICT	PIR S25754
A	81	PRO	SER	CONFLICT	PIR S25754
A	85	THR	ALA	CONFLICT	PIR S25754
A	?	-	ASN	DELETION	PIR S25754
A	97	ASP	GLY	CONFLICT	PIR S25754
A	98	VAL	ARG	CONFLICT	PIR S25754
A	99	ALA	TYR	CONFLICT	PIR S25754
B	19	THR	ILE	CONFLICT	PIR S25754
B	31	GLU	GLY	CONFLICT	PIR S25754
B	33	SER	THR	CONFLICT	PIR S25754
B	35	THR	ASN	CONFLICT	PIR S25754
B	39	HIS	GLN	CONFLICT	PIR S25754

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Chain	Residue	Modelled	Actual	Comment	Reference
B	41	SER	PRO	CONFLICT	PIR S25754
B	43	THR	ARG	CONFLICT	PIR S25754
B	50	TYR	HIS	CONFLICT	PIR S25754
B	51	GLU	SER	CONFLICT	PIR S25754
B	52	ASP	ASN	CONFLICT	PIR S25754
B	54	SER	GLN	CONFLICT	PIR S25754
B	56	ALA	PRO	CONFLICT	PIR S25754
B	60	SER	PRO	CONFLICT	PIR S25754
B	65	ALA	GLY	CONFLICT	PIR S25754
B	81	PRO	SER	CONFLICT	PIR S25754
B	85	THR	ALA	CONFLICT	PIR S25754
B	?	-	ASN	DELETION	PIR S25754
B	97	ASP	GLY	CONFLICT	PIR S25754
B	98	VAL	ARG	CONFLICT	PIR S25754
B	99	ALA	TYR	CONFLICT	PIR S25754

- Molecule 2 is water.

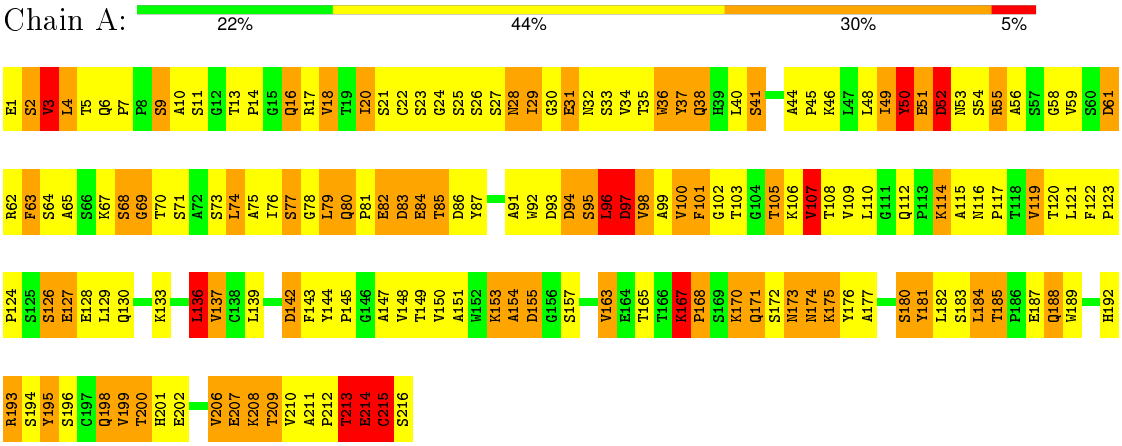
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	112	Total O 112 112	0	0
2	B	152	Total O 152 152	0	0

3 Residue-property plots (i)

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

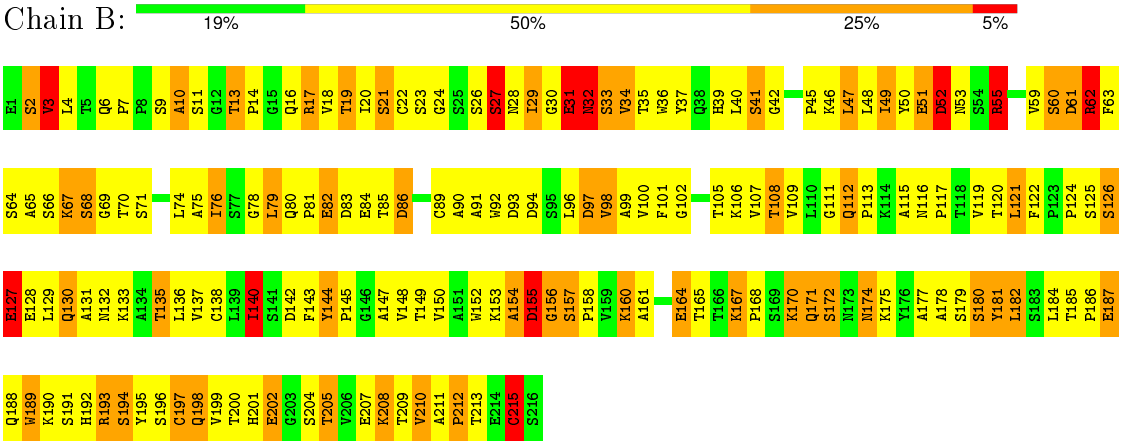
- Molecule 1: LOC - LAMBDA 1 TYPE LIGHT-CHAIN DIMER

Chain A:



- Molecule 1: LOC - LAMBDA 1 TYPE LIGHT-CHAIN DIMER

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	118.92Å 73.55Å 49.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 9.99 – 2.25	Depositor EDS
% Data completeness (in resolution range)	60.0 (10.00-2.40) 49.2 (9.99-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$	-	Xtriage
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.155 , (Not available) 0.151 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	26.4	Xtriage
Anisotropy	0.558	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.13 , 101.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[1]	$< \|L\| > = 0.41$, $< L^2 > = 0.23$	Xtriage
Outliers	0 of 10327 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3460	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	1.02	0/1628	1.99	47/2226 (2.1%)
1	B	1.02	0/1628	2.01	55/2226 (2.5%)
All	All	1.02	0/3256	2.00	102/4452 (2.3%)

There are no bond length outliers.

All (102) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	127	GLU	CA-CB-CG	15.46	147.42	113.40
1	B	193	ARG	NE-CZ-NH1	13.99	127.29	120.30
1	B	97	ASP	CB-CG-OD1	-10.22	109.10	118.30
1	A	55	ARG	NE-CZ-NH1	-9.62	115.49	120.30
1	B	55	ARG	NE-CZ-NH2	-9.42	115.59	120.30
1	B	97	ASP	CB-CG-OD2	8.54	125.99	118.30
1	A	142	ASP	CB-CG-OD2	-8.10	111.01	118.30
1	A	50	TYR	CA-CB-CG	7.73	128.08	113.40
1	A	193	ARG	CD-NE-CZ	7.60	134.24	123.60
1	B	82	GLU	OE1-CD-OE2	7.47	132.27	123.30
1	B	152	TRP	CA-CB-CG	7.40	127.75	113.70
1	A	69	GLY	N-CA-C	-7.29	94.87	113.10
1	B	79	LEU	CA-CB-CG	7.21	131.88	115.30
1	A	55	ARG	NE-CZ-NH2	7.15	123.88	120.30
1	B	160	LYS	N-CA-CB	7.02	123.23	110.60
1	A	163	VAL	CB-CA-C	6.97	124.65	111.40
1	A	52	ASP	CB-CG-OD2	-6.91	112.08	118.30
1	A	153	LYS	CA-CB-CG	6.91	128.60	113.40
1	A	97	ASP	N-CA-C	-6.87	92.44	111.00
1	A	94	ASP	CB-CG-OD2	-6.84	112.14	118.30
1	A	175	LYS	CB-CA-C	6.73	123.86	110.40
1	B	131	ALA	O-C-N	6.66	133.35	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	181	TYR	CB-CG-CD2	6.64	124.99	121.00
1	A	52	ASP	CB-CG-OD1	6.51	124.16	118.30
1	B	75	ALA	O-C-N	6.37	132.88	122.70
1	A	95	SER	N-CA-CB	6.29	119.93	110.50
1	B	51	GLU	CA-CB-CG	6.25	127.15	113.40
1	A	184	LEU	CA-CB-CG	6.24	129.66	115.30
1	B	112	GLN	CB-CG-CD	6.24	127.83	111.60
1	A	193	ARG	NE-CZ-NH1	6.09	123.35	120.30
1	A	142	ASP	N-CA-C	6.08	127.41	111.00
1	B	155	ASP	N-CA-CB	6.08	121.53	110.60
1	A	127	GLU	N-CA-CB	5.94	121.30	110.60
1	B	215	CYS	N-CA-CB	5.93	121.27	110.60
1	A	196	SER	O-C-N	5.92	132.16	122.70
1	B	62	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	B	178	ALA	CB-CA-C	5.82	118.83	110.10
1	B	180	SER	O-C-N	5.79	131.97	122.70
1	A	176	TYR	O-C-N	5.79	131.96	122.70
1	B	171	GLN	CA-CB-CG	5.78	126.11	113.40
1	B	68	SER	N-CA-CB	5.72	119.07	110.50
1	A	136	LEU	O-C-N	5.71	131.84	122.70
1	A	188	GLN	N-CA-CB	5.71	120.87	110.60
1	B	144	TYR	CA-CB-CG	-5.69	102.59	113.40
1	B	155	ASP	CB-CG-OD1	-5.64	113.23	118.30
1	B	198	GLN	CA-CB-CG	5.63	125.80	113.40
1	B	32	ASN	CB-CA-C	5.63	121.66	110.40
1	B	97	ASP	N-CA-CB	-5.62	100.48	110.60
1	B	2	SER	C-N-CA	5.62	135.74	121.70
1	A	167	LYS	CA-CB-CG	5.59	125.69	113.40
1	B	179	SER	CB-CA-C	5.59	120.72	110.10
1	B	27	SER	N-CA-CB	-5.59	102.12	110.50
1	B	140	ILE	O-C-N	5.58	131.64	122.70
1	B	86	ASP	CB-CG-OD1	5.55	123.29	118.30
1	B	121	LEU	C-N-CA	5.55	135.57	121.70
1	A	84	GLU	CG-CD-OE1	-5.54	107.21	118.30
1	B	64	SER	N-CA-CB	-5.53	102.21	110.50
1	B	61	ASP	O-C-N	5.52	131.54	122.70
1	A	37	TYR	CB-CG-CD2	-5.52	117.69	121.00
1	B	62	ARG	NE-CZ-NH1	5.51	123.05	120.30
1	A	154	ALA	O-C-N	5.50	131.50	122.70
1	A	107	VAL	O-C-N	5.49	131.49	122.70
1	B	208	LYS	O-C-N	5.45	131.41	122.70
1	A	61	ASP	CB-CG-OD2	-5.42	113.42	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	181	TYR	O-C-N	5.41	131.36	122.70
1	B	194	SER	C-N-CA	5.40	135.21	121.70
1	A	120	THR	O-C-N	5.37	131.29	122.70
1	B	142	ASP	CB-CG-OD1	5.35	123.11	118.30
1	B	98	VAL	O-C-N	5.33	131.24	122.70
1	B	156	GLY	O-C-N	5.33	131.23	122.70
1	B	31	GLU	CG-CD-OE2	5.33	128.95	118.30
1	B	155	ASP	CA-C-N	-5.29	105.63	116.20
1	A	83	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	A	36	TRP	CA-CB-CG	5.24	123.66	113.70
1	B	181	TYR	N-CA-CB	5.24	120.02	110.60
1	B	31	GLU	CG-CD-OE1	-5.23	107.83	118.30
1	A	50	TYR	CB-CA-C	5.21	120.81	110.40
1	A	185	THR	O-C-N	5.20	130.98	121.10
1	B	154	ALA	CB-CA-C	5.20	117.89	110.10
1	A	83	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	184	LEU	CA-C-N	-5.19	105.78	117.20
1	A	63	PHE	N-CA-CB	5.17	119.91	110.60
1	A	171	GLN	CB-CA-C	5.17	120.74	110.40
1	A	176	TYR	CA-CB-CG	-5.17	103.57	113.40
1	A	115	ALA	C-N-CA	5.16	134.59	121.70
1	B	125	SER	CA-CB-OG	5.15	125.09	111.20
1	A	198	GLN	OE1-CD-NE2	5.13	133.71	121.90
1	B	55	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	B	193	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	B	47	LEU	CB-CA-C	5.10	119.89	110.20
1	A	80	GLN	N-CA-C	-5.09	97.25	111.00
1	B	138	CYS	N-CA-C	-5.06	97.34	111.00
1	B	82	GLU	CG-CD-OE2	-5.05	108.19	118.30
1	A	85	THR	CA-CB-CG2	5.05	119.47	112.40
1	B	138	CYS	O-C-N	5.05	130.78	122.70
1	A	58	GLY	N-CA-C	-5.04	100.49	113.10
1	B	157	SER	O-C-N	5.04	130.68	121.10
1	B	155	ASP	CB-CA-C	-5.04	100.32	110.40
1	A	206	VAL	O-C-N	5.02	130.74	122.70
1	B	170	LYS	O-C-N	5.02	130.74	122.70
1	B	209	THR	O-C-N	5.02	130.74	122.70
1	A	193	ARG	NH1-CZ-NH2	-5.02	113.88	119.40

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	1598	0	1536	251	0
1	B	1598	0	1537	219	0
2	A	112	0	0	15	0
2	B	152	0	0	26	0
All	All	3460	0	3073	454	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 72.

All (454) 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:A:38:GLN:HG3	1:A:85:THR:CG2	1.62	1.27
1:B:52:ASP:HB2	2:B:742:HOH:O	1.28	1.26
1:B:84:GLU:OE1	1:B:170:LYS:HE2	1.46	1.12
1:A:38:GLN:CG	1:A:85:THR:HG21	1.80	1.12
1:B:130:GLN:HA	1:B:130:GLN:HE21	1.00	1.11
1:A:170:LYS:HD3	1:A:170:LYS:H	1.18	1.08
1:A:13:THR:O	1:A:16:GLN:HB2	1.54	1.06
1:B:115:ALA:HB2	1:B:175:LYS:HE2	1.40	1.03
1:B:60:SER:HB2	1:B:62:ARG:HB2	1.41	1.02
1:A:74:LEU:HD13	1:A:76:ILE:HD11	1.44	0.99
1:A:170:LYS:CD	1:A:170:LYS:H	1.77	0.97
1:B:130:GLN:HA	1:B:130:GLN:NE2	1.75	0.97
1:A:24:GLY:H	1:A:29:ILE:HD11	1.29	0.96
1:A:38:GLN:HG3	1:A:85:THR:HG21	0.97	0.95
1:A:110:LEU:HD21	1:A:114:LYS:HB2	1.49	0.95
1:B:49:ILE:HD12	1:B:55:ARG:HA	1.48	0.94
1:B:136:LEU:HD13	1:B:182:LEU:HD23	1.50	0.94
1:B:62:ARG:NH2	1:B:76:ILE:HD11	1.83	0.92
1:B:46:LYS:NZ	1:B:59:VAL:HG22	1.86	0.91
1:A:148:VAL:HG23	1:A:199:VAL:HG23	1.52	0.90
1:B:119:VAL:HG21	1:B:199:VAL:HG11	1.54	0.90
1:A:99:ALA:HB3	2:A:544:HOH:O	1.69	0.90
1:B:121:LEU:CD2	1:B:210:VAL:HG22	2.03	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:GLN:CA	1:B:130:GLN:HE21	1.85	0.88
1:A:3:VAL:HA	1:A:100:VAL:HG12	1.54	0.88
1:B:93:ASP:HB2	1:B:100:VAL:CG2	2.03	0.88
1:A:153:LYS:NZ	1:A:207:GLU:OE2	2.05	0.88
1:A:36:TRP:O	1:A:48:LEU:HB2	1.73	0.88
1:A:7:PRO:O	1:A:105:THR:HB	1.74	0.87
1:A:79:LEU:HD22	1:A:109:VAL:HG22	1.55	0.87
1:A:6:GLN:HG2	1:A:105:THR:CG2	2.04	0.87
1:A:2:SER:O	1:A:3:VAL:HB	1.74	0.86
1:A:29:ILE:O	1:A:67:LYS:NZ	2.08	0.86
1:B:115:ALA:CB	1:B:175:LYS:HE2	2.06	0.85
1:B:37:TYR:CE1	1:B:47:LEU:HD12	2.12	0.85
1:B:184:LEU:HD22	1:B:188:GLN:HB3	1.57	0.84
1:A:170:LYS:HD3	1:A:170:LYS:N	1.90	0.84
1:B:86:ASP:OD1	1:B:106:LYS:HG3	1.78	0.84
1:A:92:TRP:HZ2	1:A:97:ASP:O	1.59	0.84
1:A:24:GLY:H	1:A:29:ILE:CD1	1.91	0.83
1:B:3:VAL:HG12	1:B:101:PHE:O	1.78	0.83
1:A:93:ASP:N	1:A:98:VAL:O	2.10	0.83
1:A:30:GLY:O	1:A:31:GLU:HB2	1.78	0.83
1:A:92:TRP:CZ2	1:A:97:ASP:O	2.32	0.82
1:A:3:VAL:N	1:A:100:VAL:HG11	1.96	0.81
1:B:19:THR:HA	1:B:74:LEU:O	1.81	0.81
1:B:60:SER:C	1:B:62:ARG:H	1.80	0.81
1:A:16:GLN:O	1:A:79:LEU:HB2	1.81	0.80
1:B:4:LEU:O	1:B:102:GLY:HA2	1.82	0.80
1:B:119:VAL:HG21	1:B:199:VAL:CG1	2.12	0.80
1:B:11:SER:HB3	1:B:108:THR:HG23	1.62	0.79
1:A:74:LEU:HD13	1:A:76:ILE:CD1	2.12	0.79
1:A:189:TRP:CZ2	1:A:212:PRO:HA	2.18	0.79
1:A:24:GLY:N	1:A:29:ILE:HD11	1.98	0.78
1:B:84:GLU:OE1	1:B:170:LYS:CE	2.28	0.78
1:B:124:PRO:HD3	1:B:136:LEU:HG	1.65	0.78
1:B:119:VAL:CG2	1:B:199:VAL:HG11	2.15	0.77
1:B:10:ALA:O	1:B:107:VAL:HA	1.85	0.76
1:A:128:GLU:OE2	1:A:133:LYS:HE3	1.85	0.76
1:A:20:ILE:CG2	1:A:105:THR:HG21	2.15	0.76
1:A:38:GLN:HG3	1:A:85:THR:HG23	1.67	0.76
1:B:46:LYS:HZ1	1:B:59:VAL:HG22	1.51	0.76
1:A:6:GLN:NE2	1:A:87:TYR:O	2.20	0.75
1:A:148:VAL:HG22	1:A:149:THR:N	2.00	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ARG:NH2	1:A:83:ASP:OD1	2.18	0.75
1:A:96:LEU:CB	1:A:98:VAL:HG22	2.17	0.75
1:B:84:GLU:HG2	1:B:109:VAL:H	1.50	0.75
1:B:126:SER:O	1:B:128:GLU:N	2.20	0.75
1:A:168:PRO:HD3	2:A:618:HOH:O	1.87	0.75
1:B:121:LEU:O	2:B:605:HOH:O	2.06	0.74
1:B:192:HIS:CD2	1:B:195:TYR:OH	2.41	0.74
1:A:96:LEU:HB3	1:A:98:VAL:HG22	1.69	0.74
1:B:121:LEU:HG	1:B:210:VAL:CG2	2.17	0.74
1:A:149:THR:HB	1:A:200:THR:HG23	1.70	0.73
1:A:63:PHE:CD2	1:A:76:ILE:HG13	2.23	0.73
1:B:201:HIS:O	1:B:202:GLU:HB2	1.87	0.73
1:B:126:SER:O	1:B:127:GLU:C	2.27	0.72
1:A:119:VAL:HG11	1:A:199:VAL:CG1	2.19	0.72
1:A:148:VAL:CG2	1:A:199:VAL:HG23	2.18	0.72
1:A:119:VAL:HG11	1:A:199:VAL:HG13	1.71	0.72
1:B:170:LYS:HG2	2:B:828:HOH:O	1.88	0.72
1:B:192:HIS:HD2	1:B:195:TYR:OH	1.73	0.72
1:B:62:ARG:HH22	1:B:83:ASP:CG	1.94	0.72
1:B:13:THR:HG21	2:B:764:HOH:O	1.90	0.72
1:B:126:SER:O	1:B:129:LEU:N	2.22	0.71
1:A:148:VAL:CG2	1:A:149:THR:N	2.52	0.71
1:B:160:LYS:HE3	2:B:543:HOH:O	1.91	0.71
1:A:181:TYR:OH	1:B:171:GLN:OE1	2.09	0.70
1:B:192:HIS:NE2	2:B:849:HOH:O	2.24	0.70
1:B:208:LYS:HD3	2:B:808:HOH:O	1.91	0.70
1:B:205:THR:CG2	2:B:874:HOH:O	2.39	0.69
1:A:51:GLU:O	1:A:52:ASP:HB2	1.92	0.69
1:A:25:SER:N	1:A:28:ASN:OD1	2.25	0.69
1:B:121:LEU:HG	1:B:210:VAL:HG22	1.73	0.68
1:B:147:ALA:O	1:B:148:VAL:HG13	1.92	0.68
1:B:136:LEU:HD11	1:B:184:LEU:HD12	1.74	0.68
1:B:65:ALA:HB2	1:B:74:LEU:HD12	1.74	0.68
1:B:93:ASP:HB2	1:B:100:VAL:HG21	1.75	0.68
1:B:121:LEU:CG	1:B:210:VAL:HG22	2.24	0.68
1:B:144:TYR:CD1	1:B:144:TYR:C	2.67	0.67
1:B:182:LEU:HD21	1:B:184:LEU:HD11	1.75	0.67
1:A:214:GLU:C	1:A:215:CYS:SG	2.73	0.67
1:B:153:LYS:HA	1:B:158:PRO:HA	1.75	0.67
1:A:7:PRO:HD2	1:A:21:SER:O	1.95	0.67
1:B:60:SER:C	1:B:62:ARG:N	2.48	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:LEU:CD1	1:B:184:LEU:HD12	2.26	0.66
1:B:35:THR:HG23	1:B:49:ILE:O	1.94	0.66
1:A:171:GLN:HB2	1:B:164:GLU:HG2	1.76	0.66
1:A:112:GLN:HE22	1:A:175:LYS:HG2	1.60	0.66
1:A:101:PHE:CD1	1:A:101:PHE:N	2.62	0.66
1:A:18:VAL:O	1:A:75:ALA:HA	1.96	0.66
1:A:84:GLU:HB3	1:A:107:VAL:O	1.95	0.66
1:A:116:ASN:ND2	2:A:602:HOH:O	2.14	0.66
1:A:6:GLN:HG2	1:A:105:THR:HG22	1.78	0.66
1:A:122:PHE:CE1	1:B:137:VAL:HG21	2.31	0.66
1:A:171:GLN:NE2	1:B:181:TYR:OH	2.28	0.65
1:A:63:PHE:HD2	1:A:76:ILE:HG13	1.60	0.65
1:A:94:ASP:C	1:A:96:LEU:H	1.99	0.65
1:A:4:LEU:O	1:A:102:GLY:HA2	1.95	0.65
1:B:37:TYR:CE1	1:B:47:LEU:CD1	2.79	0.65
1:B:52:ASP:C	2:B:742:HOH:O	2.36	0.65
1:A:41:SER:HB3	2:A:560:HOH:O	1.96	0.65
1:A:34:VAL:HG23	1:A:67:LYS:HE2	1.79	0.64
1:A:70:THR:O	1:A:70:THR:HG22	1.98	0.64
1:B:196:SER:HA	1:B:208:LYS:O	1.96	0.64
1:A:209:THR:HG23	1:A:210:VAL:N	2.13	0.64
1:B:62:ARG:NH2	1:B:83:ASP:OD2	2.30	0.64
1:B:148:VAL:O	2:B:823:HOH:O	2.15	0.64
1:A:136:LEU:HD11	1:A:184:LEU:CD1	2.28	0.63
1:B:80:GLN:O	1:B:109:VAL:HG21	1.97	0.63
1:B:205:THR:HG21	2:B:874:HOH:O	1.96	0.63
1:A:38:GLN:HB2	1:A:87:TYR:CE2	2.33	0.63
1:B:20:ILE:N	1:B:20:ILE:HD13	2.13	0.63
1:A:117:PRO:HB3	1:A:143:PHE:HB3	1.79	0.63
1:B:211:ALA:O	1:B:213:THR:N	2.32	0.63
1:A:84:GLU:HA	1:A:107:VAL:HG22	1.81	0.62
1:A:150:VAL:HG12	1:A:151:ALA:N	2.13	0.62
1:A:48:LEU:HD11	1:A:87:TYR:CD2	2.35	0.62
1:B:186:PRO:O	1:B:190:LYS:HG2	2.00	0.62
1:A:2:SER:C	1:A:100:VAL:HG11	2.20	0.62
1:B:121:LEU:HD21	1:B:210:VAL:HG22	1.81	0.62
1:B:97:ASP:O	1:B:97:ASP:OD1	2.18	0.61
1:A:20:ILE:CG2	1:A:105:THR:CG2	2.78	0.61
1:A:24:GLY:O	1:A:70:THR:HG22	2.01	0.61
1:B:184:LEU:CD2	1:B:188:GLN:HB3	2.28	0.61
1:A:150:VAL:HG12	1:A:151:ALA:H	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:SER:HB3	2:B:720:HOH:O	1.98	0.61
1:A:36:TRP:O	1:A:48:LEU:N	2.29	0.61
1:A:6:GLN:OE1	1:A:102:GLY:HA3	2.00	0.61
1:B:130:GLN:O	2:B:842:HOH:O	2.16	0.61
1:A:2:SER:HB3	1:A:100:VAL:HG11	1.81	0.61
1:A:182:LEU:HD11	2:A:614:HOH:O	2.01	0.61
1:A:62:ARG:HH22	1:A:83:ASP:CG	2.03	0.61
1:B:113:PRO:O	2:B:804:HOH:O	2.16	0.61
1:A:76:ILE:HG22	1:A:76:ILE:O	2.00	0.61
1:B:46:LYS:HZ3	1:B:59:VAL:HG22	1.63	0.61
1:A:92:TRP:CE3	1:A:99:ALA:HB2	2.36	0.61
1:B:19:THR:C	1:B:20:ILE:HD13	2.21	0.61
1:A:22:CYS:O	1:A:29:ILE:CD1	2.47	0.60
1:B:48:LEU:HD23	1:B:74:LEU:HD21	1.82	0.60
1:B:189:TRP:CZ2	1:B:212:PRO:HA	2.36	0.60
1:A:209:THR:CG2	1:A:210:VAL:N	2.65	0.60
1:A:6:GLN:HG2	1:A:105:THR:HG21	1.84	0.60
1:B:4:LEU:CD2	1:B:29:ILE:HD11	2.31	0.60
1:B:32:ASN:ND2	1:B:92:TRP:O	2.35	0.60
1:A:92:TRP:CZ2	1:A:97:ASP:C	2.75	0.60
1:A:121:LEU:HD23	1:A:209:THR:HA	1.84	0.60
1:A:119:VAL:CG2	1:A:208:LYS:HB2	2.32	0.59
1:A:74:LEU:CD1	1:A:76:ILE:HD11	2.25	0.59
1:B:187:GLU:O	1:B:188:GLN:C	2.40	0.59
1:B:24:GLY:O	1:B:70:THR:HB	2.03	0.59
1:B:52:ASP:CB	2:B:742:HOH:O	2.10	0.59
1:B:90:ALA:HA	1:B:100:VAL:O	2.02	0.59
1:A:184:LEU:HD22	1:A:189:TRP:HA	1.84	0.59
1:A:185:THR:OG1	1:A:188:GLN:HG3	2.02	0.59
1:A:121:LEU:HD23	1:A:209:THR:CA	2.33	0.59
1:A:3:VAL:CA	1:A:100:VAL:HG12	2.29	0.59
1:B:4:LEU:HD21	1:B:29:ILE:HG13	1.85	0.58
1:B:145:PRO:HA	2:B:817:HOH:O	2.03	0.58
1:A:182:LEU:HD21	1:A:195:TYR:HE1	1.68	0.58
1:A:11:SER:OG	1:A:110:LEU:HD11	2.03	0.58
1:A:148:VAL:CG2	1:A:149:THR:H	2.15	0.58
1:B:18:VAL:HG22	1:B:76:ILE:CG2	2.32	0.58
1:B:191:SER:O	1:B:191:SER:OG	2.21	0.58
1:A:119:VAL:HG22	1:A:208:LYS:HD3	1.86	0.58
1:A:122:PHE:CZ	1:B:137:VAL:CG2	2.87	0.58
1:A:80:GLN:HG2	1:A:81:PRO:HD2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ARG:NH1	2:A:551:HOH:O	2.30	0.57
1:B:96:LEU:O	1:B:97:ASP:C	2.41	0.57
1:B:182:LEU:CD2	1:B:184:LEU:HD11	2.34	0.57
1:A:119:VAL:HA	1:A:139:LEU:O	2.05	0.57
1:A:49:ILE:HG21	1:A:65:ALA:HB3	1.86	0.57
1:A:97:ASP:H	1:A:98:VAL:HG22	1.69	0.57
1:A:126:SER:O	1:A:127:GLU:C	2.40	0.57
1:A:5:THR:O	1:A:22:CYS:HA	2.05	0.57
1:A:184:LEU:HD22	1:A:189:TRP:CA	2.33	0.57
1:A:144:TYR:HA	1:A:145:PRO:C	2.24	0.57
1:A:45:PRO:HG2	1:B:101:PHE:CD2	2.40	0.56
1:A:93:ASP:HB2	1:A:100:VAL:HG23	1.87	0.56
1:A:82:GLU:HG3	2:A:555:HOH:O	2.05	0.56
1:B:51:GLU:O	1:B:53:ASN:N	2.31	0.56
1:B:35:THR:HG21	1:B:47:LEU:HD11	1.86	0.56
1:B:121:LEU:HG	1:B:210:VAL:HG21	1.88	0.56
1:A:25:SER:C	1:A:27:SER:H	2.08	0.56
1:A:29:ILE:HD12	1:A:71:SER:HA	1.88	0.56
1:A:24:GLY:O	1:A:70:THR:CG2	2.54	0.55
1:A:182:LEU:HD21	1:A:195:TYR:CE1	2.42	0.55
1:B:62:ARG:HH21	1:B:76:ILE:HD11	1.70	0.55
1:A:124:PRO:HB2	1:A:129:LEU:HD13	1.88	0.55
1:A:20:ILE:HD11	1:A:76:ILE:HD13	1.88	0.55
1:A:7:PRO:O	1:A:105:THR:CB	2.51	0.55
1:B:84:GLU:HG2	1:B:108:THR:HA	1.88	0.55
1:A:149:THR:O	1:A:199:VAL:HA	2.05	0.55
1:B:28:ASN:O	1:B:29:ILE:C	2.45	0.55
1:A:136:LEU:O	1:A:181:TYR:HA	2.06	0.55
1:B:49:ILE:HD11	1:B:55:ARG:HG3	1.87	0.55
1:B:80:GLN:O	1:B:81:PRO:C	2.45	0.55
1:A:6:GLN:HE21	1:A:105:THR:HG23	1.71	0.55
1:A:93:ASP:HB2	1:A:100:VAL:CG2	2.37	0.55
1:A:79:LEU:HD22	1:A:109:VAL:CG2	2.31	0.54
1:A:29:ILE:HG21	1:A:71:SER:CA	2.37	0.54
1:A:3:VAL:N	1:A:100:VAL:CG1	2.69	0.54
1:A:136:LEU:HD11	1:A:184:LEU:HD11	1.89	0.54
1:A:155:ASP:OD1	1:A:193:ARG:HB3	2.07	0.54
1:B:93:ASP:O	1:B:97:ASP:N	2.41	0.54
1:A:212:PRO:O	2:A:624:HOH:O	2.18	0.54
1:A:37:TYR:HA	1:A:46:LYS:O	2.07	0.54
1:B:50:TYR:O	1:B:51:GLU:C	2.44	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:GLU:OE1	1:B:133:LYS:HE3	2.06	0.54
1:A:1:PCA:O	1:A:2:SER:CB	2.55	0.54
1:B:201:HIS:O	1:B:202:GLU:CB	2.54	0.54
1:B:33:SER:HA	1:B:52:ASP:OD1	2.07	0.53
1:B:49:ILE:CD1	1:B:55:ARG:HG3	2.38	0.53
1:B:200:THR:OG1	1:B:205:THR:OG1	2.20	0.53
1:B:155:ASP:OD1	1:B:193:ARG:HB2	2.08	0.53
1:B:76:ILE:HD12	1:B:76:ILE:C	2.29	0.53
1:A:3:VAL:CA	1:A:100:VAL:CG1	2.86	0.53
1:A:3:VAL:HA	1:A:100:VAL:CG1	2.32	0.53
1:A:121:LEU:CD2	1:A:209:THR:HA	2.39	0.53
1:B:37:TYR:CD1	1:B:47:LEU:HA	2.44	0.53
1:B:18:VAL:HG22	1:B:76:ILE:HG23	1.91	0.53
1:A:62:ARG:O	1:A:77:SER:HB2	2.08	0.53
1:A:172:SER:C	1:A:174:ASN:H	2.12	0.53
1:A:184:LEU:CD2	1:A:189:TRP:HA	2.38	0.53
1:B:4:LEU:HD23	1:B:29:ILE:HD11	1.91	0.53
1:A:193:ARG:HB3	2:A:613:HOH:O	2.08	0.53
1:A:96:LEU:O	1:A:97:ASP:HB2	2.08	0.52
1:A:34:VAL:O	1:A:52:ASP:N	2.35	0.52
1:B:130:GLN:HB2	2:B:867:HOH:O	2.09	0.52
1:B:121:LEU:CG	1:B:210:VAL:CG2	2.84	0.52
1:B:22:CYS:O	1:B:29:ILE:CD1	2.57	0.52
1:B:92:TRP:HA	1:B:99:ALA:HA	1.91	0.52
1:A:208:LYS:HE3	2:A:628:HOH:O	2.08	0.52
1:A:44:ALA:CB	1:B:102:GLY:O	2.57	0.52
1:B:208:LYS:CD	2:B:808:HOH:O	2.53	0.52
1:B:121:LEU:HD21	1:B:210:VAL:CG2	2.39	0.52
1:A:215:CYS:O	1:A:216:SER:HB2	2.09	0.52
1:A:85:THR:HG22	1:A:86:ASP:O	2.10	0.51
1:B:92:TRP:HZ2	1:B:97:ASP:OD1	1.94	0.51
1:A:20:ILE:HG22	1:A:105:THR:HG21	1.91	0.51
1:B:187:GLU:O	1:B:191:SER:HB3	2.11	0.51
1:B:188:GLN:O	1:B:192:HIS:HD2	1.93	0.51
1:B:29:ILE:HG12	1:B:34:VAL:HG23	1.92	0.51
1:A:136:LEU:HD12	1:A:136:LEU:N	2.26	0.51
1:A:6:GLN:NE2	1:A:105:THR:HG23	2.26	0.51
1:A:11:SER:HB2	1:A:110:LEU:HD12	1.92	0.51
1:A:48:LEU:HD11	1:A:87:TYR:HD2	1.76	0.50
1:B:60:SER:HB2	1:B:62:ARG:CB	2.27	0.50
1:A:84:GLU:HA	1:A:107:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:LYS:HA	1:B:71:SER:O	2.12	0.50
1:B:154:ALA:O	1:B:156:GLY:N	2.44	0.50
1:B:37:TYR:HA	1:B:46:LYS:O	2.11	0.50
1:B:132:ASN:HA	1:B:186:PRO:HG2	1.94	0.50
1:B:7:PRO:O	1:B:105:THR:OG1	2.27	0.50
1:A:20:ILE:HD11	1:A:76:ILE:CD1	2.41	0.50
1:A:85:THR:CG2	1:A:86:ASP:H	2.24	0.50
1:A:85:THR:CG2	1:A:86:ASP:N	2.75	0.50
1:A:126:SER:O	1:A:130:GLN:HG3	2.12	0.50
1:B:126:SER:C	1:B:128:GLU:N	2.64	0.50
1:A:92:TRP:CD1	1:A:93:ASP:O	2.66	0.49
1:A:6:GLN:CG	1:A:105:THR:HG22	2.42	0.49
1:B:182:LEU:HD21	1:B:184:LEU:CD1	2.41	0.49
1:A:48:LEU:HD11	1:A:87:TYR:CE2	2.48	0.49
1:A:82:GLU:HB2	2:A:542:HOH:O	2.11	0.49
1:B:93:ASP:HB2	1:B:100:VAL:HG23	1.92	0.49
1:B:144:TYR:CD1	1:B:145:PRO:N	2.81	0.49
1:B:35:THR:HG21	1:B:47:LEU:CD1	2.43	0.49
1:A:14:PRO:HG3	1:A:109:VAL:CG1	2.43	0.49
1:B:92:TRP:HZ3	2:B:771:HOH:O	1.95	0.49
1:A:33:SER:OG	1:A:51:GLU:OE1	2.31	0.49
1:A:44:ALA:HB2	1:B:102:GLY:O	2.13	0.49
1:B:140:ILE:HG22	1:B:143:PHE:CD2	2.48	0.49
1:B:135:THR:CG2	1:B:181:TYR:HD2	2.26	0.48
1:A:17:ARG:NH1	2:A:523:HOH:O	2.46	0.48
1:A:168:PRO:HA	1:A:177:ALA:O	2.14	0.48
1:B:97:ASP:OD1	1:B:97:ASP:C	2.51	0.48
1:A:182:LEU:CD2	1:A:195:TYR:CE1	2.96	0.48
1:B:40:LEU:HD13	1:B:85:THR:HG22	1.96	0.48
1:A:214:GLU:O	1:A:215:CYS:C	2.52	0.48
1:B:150:VAL:CG2	1:B:197:CYS:SG	3.01	0.48
1:B:196:SER:CA	1:B:208:LYS:O	2.62	0.48
1:A:165:THR:HG23	1:A:180:SER:HB2	1.94	0.48
1:A:48:LEU:CD1	1:A:87:TYR:HD2	2.26	0.48
1:A:81:PRO:HA	1:A:109:VAL:HG21	1.96	0.48
1:B:34:VAL:HG23	1:B:91:ALA:HB2	1.95	0.48
1:B:14:PRO:HD3	1:B:111:GLY:H	1.79	0.48
1:B:136:LEU:O	1:B:181:TYR:HA	2.14	0.48
1:A:119:VAL:HG11	1:A:199:VAL:HG11	1.96	0.48
1:B:96:LEU:HB3	1:B:98:VAL:HG22	1.95	0.48
1:B:197:CYS:O	1:B:207:GLU:HA	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ALA:O	1:B:148:VAL:CG1	2.60	0.47
1:A:38:GLN:CG	1:A:85:THR:CG2	2.57	0.47
1:B:4:LEU:HD12	1:B:100:VAL:HB	1.97	0.47
1:B:2:SER:OG	1:B:3:VAL:N	2.47	0.47
1:B:175:LYS:HD3	2:B:844:HOH:O	2.13	0.47
1:B:172:SER:C	1:B:174:ASN:H	2.16	0.47
1:B:192:HIS:HD2	1:B:195:TYR:HH	1.61	0.47
1:B:4:LEU:HD22	1:B:22:CYS:SG	2.55	0.47
1:B:9:SER:HA	1:B:106:LYS:O	2.15	0.47
1:A:49:ILE:N	1:A:49:ILE:CD1	2.78	0.47
1:B:165:THR:OG1	1:B:180:SER:OG	2.11	0.47
1:B:132:ASN:OD1	1:B:186:PRO:HD2	2.14	0.47
1:A:25:SER:O	1:A:27:SER:N	2.36	0.47
1:B:51:GLU:C	1:B:53:ASN:H	2.16	0.47
1:A:155:ASP:OD1	1:A:193:ARG:CB	2.63	0.46
1:B:182:LEU:CD2	1:B:184:LEU:CD1	2.93	0.46
1:B:119:VAL:O	1:B:208:LYS:HE3	2.15	0.46
1:A:45:PRO:HG2	1:B:101:PHE:CG	2.50	0.46
1:A:10:ALA:O	1:A:107:VAL:HA	2.15	0.46
1:A:155:ASP:OD2	1:A:192:HIS:HD2	1.98	0.46
1:B:35:THR:HG22	1:B:36:TRP:N	2.31	0.46
1:A:92:TRP:HA	1:A:99:ALA:HA	1.98	0.46
1:A:112:GLN:HE22	1:A:175:LYS:CG	2.26	0.46
1:A:122:PHE:CZ	1:B:137:VAL:HG21	2.50	0.46
1:A:99:ALA:CB	2:A:544:HOH:O	2.44	0.46
1:B:32:ASN:N	1:B:32:ASN:ND2	2.61	0.46
1:A:93:ASP:CB	1:A:98:VAL:O	2.63	0.46
1:B:92:TRP:CZ2	1:B:97:ASP:OD1	2.69	0.46
1:A:182:LEU:CD2	1:A:195:TYR:HE1	2.29	0.46
1:A:93:ASP:CA	1:A:98:VAL:O	2.64	0.46
1:A:213:THR:HG21	1:B:215:CYS:SG	2.56	0.46
1:B:124:PRO:HB2	1:B:129:LEU:HD13	1.98	0.46
1:A:150:VAL:CG1	1:A:151:ALA:N	2.79	0.46
1:A:142:ASP:OD1	1:A:171:GLN:OE1	2.34	0.45
1:B:6:GLN:OE1	1:B:89:CYS:SG	2.74	0.45
1:A:36:TRP:C	1:A:48:LEU:HB2	2.34	0.45
1:A:11:SER:OG	1:A:110:LEU:CD1	2.65	0.45
1:A:68:SER:C	1:A:69:GLY:O	2.51	0.45
1:B:187:GLU:O	1:B:189:TRP:N	2.49	0.45
1:B:121:LEU:CD2	1:B:210:VAL:CG2	2.83	0.45
1:B:65:ALA:HB2	1:B:74:LEU:CD1	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:TYR:CZ	1:B:47:LEU:HD12	2.51	0.45
1:B:79:LEU:HD13	1:B:109:VAL:HG22	1.99	0.45
1:A:92:TRP:HD1	1:A:93:ASP:O	2.00	0.45
1:A:150:VAL:HA	1:A:198:GLN:O	2.16	0.45
1:B:53:ASN:ND2	2:B:742:HOH:O	2.49	0.45
1:A:5:THR:O	1:A:23:SER:N	2.48	0.45
1:B:62:ARG:CZ	1:B:76:ILE:HD11	2.46	0.45
1:A:20:ILE:CD1	1:A:74:LEU:HB3	2.48	0.44
1:B:49:ILE:CD1	1:B:55:ARG:HA	2.35	0.44
1:A:213:THR:HG23	1:A:214:GLU:O	2.16	0.44
1:A:36:TRP:O	1:A:48:LEU:CB	2.57	0.44
1:A:65:ALA:HA	1:A:73:SER:O	2.17	0.44
1:A:170:LYS:HD2	1:A:170:LYS:H	1.75	0.44
1:A:80:GLN:HG2	1:A:81:PRO:CD	2.46	0.44
1:A:153:LYS:HA	1:A:157:SER:O	2.17	0.44
1:B:6:GLN:HA	1:B:21:SER:O	2.18	0.44
1:A:9:SER:HB2	1:A:147:ALA:HB3	2.00	0.44
1:A:101:PHE:HB2	1:B:45:PRO:HD2	2.00	0.44
1:A:3:VAL:HG22	1:A:101:PHE:O	2.18	0.44
1:A:209:THR:HG23	1:A:210:VAL:H	1.82	0.44
1:A:184:LEU:HB3	1:A:185:THR:H	1.61	0.44
1:A:206:VAL:CG2	2:A:601:HOH:O	2.66	0.44
1:A:48:LEU:O	1:A:59:VAL:HG21	2.18	0.43
1:A:148:VAL:HG23	1:A:149:THR:H	1.81	0.43
1:A:50:TYR:N	1:A:54:SER:O	2.49	0.43
1:A:211:ALA:HA	1:A:212:PRO:HD3	1.72	0.43
1:B:17:ARG:HB2	2:B:739:HOH:O	2.18	0.43
1:A:214:GLU:C	1:A:215:CYS:O	2.57	0.43
1:A:150:VAL:CG1	1:A:151:ALA:H	2.29	0.43
1:B:39:HIS:O	1:B:41:SER:N	2.52	0.43
1:A:14:PRO:HG3	1:A:109:VAL:HG13	2.00	0.43
1:A:24:GLY:HA3	1:A:29:ILE:CG1	2.48	0.43
1:A:32:ASN:O	1:A:67:LYS:NZ	2.48	0.43
1:B:63:PHE:HZ	2:B:718:HOH:O	2.01	0.43
1:B:128:GLU:O	1:B:133:LYS:O	2.37	0.43
1:A:76:ILE:N	1:A:76:ILE:HD12	2.34	0.43
1:A:91:ALA:O	1:A:99:ALA:HA	2.19	0.43
1:B:4:LEU:CD2	1:B:29:ILE:CD1	2.97	0.43
1:B:92:TRP:CD1	1:B:94:ASP:OD1	2.72	0.43
1:A:172:SER:OG	1:A:173:ASN:N	2.52	0.42
1:A:20:ILE:HG23	1:A:105:THR:HG21	1.98	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:VAL:HG21	1:B:122:PHE:CE2	2.54	0.42
1:B:53:ASN:N	2:B:742:HOH:O	2.51	0.42
1:A:171:GLN:HE21	1:B:164:GLU:HG2	1.84	0.42
1:A:49:ILE:N	1:A:49:ILE:HD12	2.35	0.42
1:B:40:LEU:O	1:B:42:GLY:N	2.53	0.42
1:B:82:GLU:HG3	1:B:174:ASN:OD1	2.20	0.42
1:A:172:SER:HB3	1:B:164:GLU:OE2	2.19	0.42
1:B:117:PRO:CD	1:B:201:HIS:CD2	3.02	0.42
1:B:14:PRO:HD3	1:B:111:GLY:N	2.35	0.42
1:B:84:GLU:CG	1:B:109:VAL:H	2.25	0.42
1:B:60:SER:HB3	1:B:62:ARG:HG3	2.01	0.42
1:A:151:ALA:O	1:A:198:GLN:HB3	2.20	0.42
1:A:170:LYS:N	1:A:170:LYS:CD	2.53	0.42
1:B:121:LEU:HD23	1:B:210:VAL:HG22	1.97	0.42
1:A:121:LEU:HD23	1:A:209:THR:C	2.40	0.42
1:A:85:THR:HG23	1:A:86:ASP:H	1.85	0.42
1:B:119:VAL:O	1:B:208:LYS:HD2	2.19	0.42
1:A:55:ARG:HG2	1:A:56:ALA:O	2.20	0.42
1:A:154:ALA:O	1:A:157:SER:HB2	2.19	0.42
1:A:4:LEU:N	1:A:4:LEU:CD1	2.80	0.41
1:A:56:ALA:HB3	1:A:59:VAL:HG21	2.01	0.41
1:B:50:TYR:CE2	1:B:55:ARG:O	2.73	0.41
1:A:70:THR:O	1:A:70:THR:CG2	2.65	0.41
1:B:27:SER:HB2	1:B:93:ASP:OD1	2.20	0.41
1:B:4:LEU:HD21	1:B:29:ILE:CG1	2.48	0.41
1:B:13:THR:O	1:B:16:GLN:HB3	2.20	0.41
1:B:112:GLN:HB3	1:B:113:PRO:CD	2.50	0.41
1:A:74:LEU:HD23	1:A:74:LEU:HA	1.91	0.41
1:B:49:ILE:HD12	1:B:49:ILE:HA	1.82	0.41
1:B:208:LYS:CE	2:B:808:HOH:O	2.69	0.41
1:A:184:LEU:HD21	1:A:195:TYR:CZ	2.56	0.41
1:B:174:ASN:HD22	1:B:174:ASN:HA	1.54	0.41
1:A:50:TYR:HD1	1:A:54:SER:O	2.03	0.41
1:A:148:VAL:CG2	1:A:199:VAL:CG2	2.95	0.41
1:A:136:LEU:HD12	1:A:182:LEU:O	2.21	0.41
1:A:70:THR:C	1:A:71:SER:OG	2.59	0.41
1:A:93:ASP:OD2	1:A:100:VAL:HG21	2.20	0.41
1:B:13:THR:HA	1:B:111:GLY:H	1.86	0.41
1:A:123:PRO:HG3	1:A:210:VAL:HG21	2.03	0.41
1:A:68:SER:HB3	1:A:69:GLY:H	1.57	0.41
1:B:68:SER:HB3	2:B:729:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ASN:O	1:A:67:LYS:CE	2.69	0.41
1:B:4:LEU:HD23	1:B:29:ILE:CD1	2.50	0.41
1:A:116:ASN:HA	1:A:201:HIS:HD2	1.85	0.41
1:A:49:ILE:HG21	1:A:65:ALA:CB	2.50	0.41
1:A:6:GLN:CG	1:A:105:THR:CG2	2.87	0.41
1:A:33:SER:CB	2:A:530:HOH:O	2.70	0.41
1:B:119:VAL:C	1:B:208:LYS:HD2	2.41	0.41
1:A:45:PRO:HD2	1:B:101:PHE:HB3	2.03	0.41
1:B:74:LEU:HA	1:B:74:LEU:HD12	1.97	0.41
1:B:117:PRO:HD3	1:B:201:HIS:CD2	2.56	0.41
1:A:25:SER:C	1:A:27:SER:N	2.73	0.41
1:B:67:LYS:HE3	1:B:69:GLY:O	2.21	0.41
1:A:4:LEU:HA	1:A:4:LEU:HD12	1.58	0.40
1:A:29:ILE:HG21	1:A:71:SER:C	2.41	0.40
1:B:140:ILE:O	1:B:177:ALA:HA	2.21	0.40
1:B:185:THR:O	1:B:186:PRO:C	2.59	0.40
1:B:167:LYS:HA	1:B:168:PRO:HD3	1.87	0.40
1:A:85:THR:HG22	1:A:87:TYR:CE1	2.57	0.40
1:B:185:THR:O	1:B:187:GLU:N	2.54	0.40
1:A:86:ASP:OD1	1:A:86:ASP:N	2.54	0.40
1:A:167:LYS:O	1:A:168:PRO:C	2.60	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	214/216 (99%)	174 (81%)	25 (12%)	15 (7%)	1 0
1	B	214/216 (99%)	171 (80%)	27 (13%)	16 (8%)	1 0
All	All	428/432 (99%)	345 (81%)	52 (12%)	31 (7%)	1 0

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	31	GLU
1	A	77	SER
1	A	96	LEU
1	A	97	ASP
1	A	214	GLU
1	B	3	VAL
1	B	41	SER
1	B	52	ASP
1	B	55	ARG
1	A	202	GLU
1	A	215	CYS
1	B	29	ILE
1	B	78	GLY
1	B	127	GLU
1	B	189	TRP
1	B	212	PRO
1	A	79	LEU
1	A	155	ASP
1	A	213	THR
1	B	10	ALA
1	B	30	GLY
1	B	31	GLU
1	B	155	ASP
1	A	26	SER
1	A	52	ASP
1	B	161	ALA
1	B	187	GLU
1	B	215	CYS
1	A	78	GLY
1	A	168	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	181/181 (100%)	126 (70%)	55 (30%)	0 0
1	B	181/181 (100%)	139 (77%)	42 (23%)	1 1
All	All	362/362 (100%)	265 (73%)	97 (27%)	0 0

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	3	VAL
1	A	4	LEU
1	A	9	SER
1	A	16	GLN
1	A	18	VAL
1	A	20	ILE
1	A	28	ASN
1	A	29	ILE
1	A	35	THR
1	A	38	GLN
1	A	40	LEU
1	A	41	SER
1	A	49	ILE
1	A	50	TYR
1	A	51	GLU
1	A	53	ASN
1	A	61	ASP
1	A	64	SER
1	A	68	SER
1	A	74	LEU
1	A	82	GLU
1	A	95	SER
1	A	96	LEU
1	A	98	VAL
1	A	100	VAL
1	A	101	PHE
1	A	103	THR
1	A	105	THR
1	A	106	LYS
1	A	107	VAL
1	A	108	THR
1	A	114	LYS
1	A	119	VAL
1	A	126	SER

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Mol	Chain	Res	Type
1	A	136	LEU
1	A	137	VAL
1	A	163	VAL
1	A	167	LYS
1	A	170	LYS
1	A	173	ASN
1	A	174	ASN
1	A	180	SER
1	A	183	SER
1	A	187	GLU
1	A	194	SER
1	A	195	TYR
1	A	199	VAL
1	A	200	THR
1	A	207	GLU
1	A	208	LYS
1	A	209	THR
1	A	213	THR
1	A	214	GLU
1	A	215	CYS
1	B	3	VAL
1	B	13	THR
1	B	17	ARG
1	B	19	THR
1	B	21	SER
1	B	23	SER
1	B	26	SER
1	B	27	SER
1	B	31	GLU
1	B	32	ASN
1	B	33	SER
1	B	34	VAL
1	B	49	ILE
1	B	52	ASP
1	B	60	SER
1	B	61	ASP
1	B	62	ARG
1	B	66	SER
1	B	67	LYS
1	B	76	ILE
1	B	108	THR
1	B	116	ASN

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Mol	Chain	Res	Type
1	B	120	THR
1	B	126	SER
1	B	127	GLU
1	B	130	GLN
1	B	135	THR
1	B	140	ILE
1	B	149	THR
1	B	157	SER
1	B	164	GLU
1	B	167	LYS
1	B	172	SER
1	B	174	ASN
1	B	182	LEU
1	B	194	SER
1	B	197	CYS
1	B	198	GLN
1	B	202	GLU
1	B	204	SER
1	B	205	THR
1	B	210	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	171	GLN
1	A	173	ASN
1	A	188	GLN
1	A	192	HIS
1	B	38	GLN
1	B	80	GLN
1	B	112	GLN
1	B	130	GLN
1	B	174	ASN
1	B	192	HIS
1	B	201	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	PCA	A	1	1	6,7,9	0.98	1 (16%)	7,8,12	1.66	2 (28%)
1	PCA	B	1	1	6,7,9	0.79	0	7,8,12	1.52	2 (28%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PCA	A	1	1	-	0/0/9/13	0/1/1/1
1	PCA	B	1	1	-	0/0/9/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	PCA	O-C	2.02	1.29	1.19

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	1	PCA	CB-CA-N	-2.53	99.81	105.01
1	A	1	PCA	O-C-CA	-2.30	119.35	125.44
1	B	1	PCA	CB-CA-C	2.68	116.43	112.76
1	A	1	PCA	CB-CA-C	3.28	117.25	112.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1	PCA	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	215/216 (99%)	-1.13	0	100	100	5, 22, 43, 56	0
1	B	215/216 (99%)	-1.23	0	100	100	5, 20, 32, 52	0
All	All	430/432 (99%)	-1.18	0	100	100	5, 21, 41, 56	0

There are no RSRZ outliers to report.

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
1	PCA	B	1	7/9	0.90	0.13	-	42,42,42,42	0
1	PCA	A	1	7/9	0.91	0.19	-	40,41,42,42	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.