



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:28 PM GMT

PDB ID : 1AXS
Title : MATURE OXY-COPE CATALYTIC ANTIBODY WITH HAPTEN
Authors : Mundorff, E.C.; Ulrich, H.D.; Stevens, R.C.
Deposited on : 1997-10-20
Resolution : 2.60 Å(reported)

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

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

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

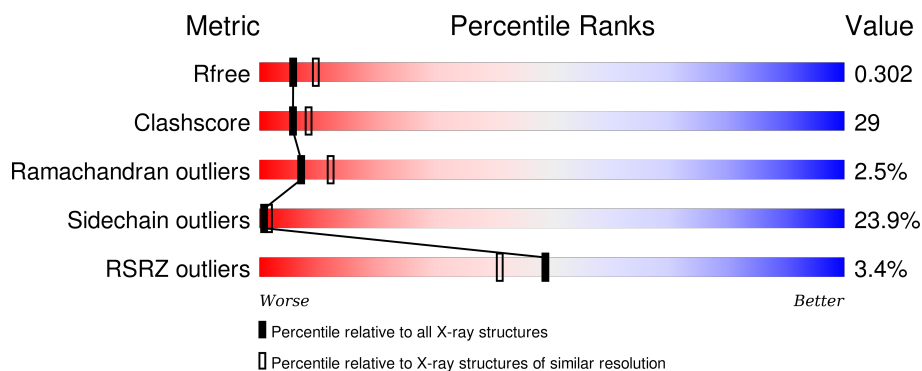
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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

Mol	Chain	Length	Quality of chain
1	A	211	<div> <div>2%</div> <div> <div>43%</div> <div>43%</div> <div>14%</div> </div> </div>
1	L	211	<div> <div>%</div> <div> <div>46%</div> <div>41%</div> <div>13%</div> </div> </div>
2	B	221	<div> <div>4%</div> <div> <div>47%</div> <div>38%</div> <div>14%</div> </div> </div>
2	H	221	<div> <div>6%</div> <div> <div>46%</div> <div>44%</div> <div>9%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	HOP	H	1006	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6887 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 OXY-COPE CATALYTIC ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1655	1041	272	336	6			
1	A	211	Total	C	N	O	S	0	0	0
			1655	1041	272	336	6			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	11	MET	LEU	CONFLICT	GB 243868
L	12	TYR	SER	CONFLICT	GB 243868
L	15	LEU	VAL	CONFLICT	GB 243868
L	17	GLU	ASP	CONFLICT	GB 243868
L	28	ASP	ASN	CONFLICT	GB 243868
L	30	ASN	ASP	CONFLICT	GB 243868
L	31	SER	LYS	CONFLICT	GB 243868
L	36	PHE	TYR	CONFLICT	GB 243868
L	43	SER	ALA	CONFLICT	GB 243868
L	46	THR	LEU	CONFLICT	GB 243868
L	50	ARG	ASN	CONFLICT	GB 243868
L	53	ARG	ASN	CONFLICT	GB 243868
L	55	VAL	GLN	CONFLICT	GB 243868
L	56	ASP	THR	CONFLICT	GB 243868
L	69	GLN	THR	CONFLICT	GB 243868
L	71	TYR	PHE	CONFLICT	GB 243868
L	72	SER	THR	CONFLICT	GB 243868
L	73	LEU	PHE	CONFLICT	GB 243868
L	79	GLU	GLN	CONFLICT	GB 243868
L	80	TYR	PRO	CONFLICT	GB 243868
L	83	MET	ILE	CONFLICT	GB 243868
L	84	GLY	ALA	CONFLICT	GB 243868
L	85	ILE	THR	CONFLICT	GB 243868
L	91	TYR	HIS	CONFLICT	GB 243868
L	92	ASP	ILE	CONFLICT	GB 243868

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Chain	Residue	Modelled	Actual	Comment	Reference
L	93	GLU	SER	CONFLICT	GB 243868
L	94	PHE	ARG	CONFLICT	GB 243868
L	96	TYR	ARG	CONFLICT	GB 243868
L	100	SER	GLN	CONFLICT	GB 243868
L	104	LEU	VAL	CONFLICT	GB 243868
A	11	MET	LEU	CONFLICT	GB 243868
A	12	TYR	SER	CONFLICT	GB 243868
A	15	LEU	VAL	CONFLICT	GB 243868
A	17	GLU	ASP	CONFLICT	GB 243868
A	28	ASP	ASN	CONFLICT	GB 243868
A	30	ASN	ASP	CONFLICT	GB 243868
A	31	SER	LYS	CONFLICT	GB 243868
A	36	PHE	TYR	CONFLICT	GB 243868
A	43	SER	ALA	CONFLICT	GB 243868
A	46	THR	LEU	CONFLICT	GB 243868
A	50	ARG	ASN	CONFLICT	GB 243868
A	53	ARG	ASN	CONFLICT	GB 243868
A	55	VAL	GLN	CONFLICT	GB 243868
A	56	ASP	THR	CONFLICT	GB 243868
A	69	GLN	THR	CONFLICT	GB 243868
A	71	TYR	PHE	CONFLICT	GB 243868
A	72	SER	THR	CONFLICT	GB 243868
A	73	LEU	PHE	CONFLICT	GB 243868
A	79	GLU	GLN	CONFLICT	GB 243868
A	80	TYR	PRO	CONFLICT	GB 243868
A	83	MET	ILE	CONFLICT	GB 243868
A	84	GLY	ALA	CONFLICT	GB 243868
A	85	ILE	THR	CONFLICT	GB 243868
A	91	TYR	HIS	CONFLICT	GB 243868
A	92	ASP	ILE	CONFLICT	GB 243868
A	93	GLU	SER	CONFLICT	GB 243868
A	94	PHE	ARG	CONFLICT	GB 243868
A	96	TYR	ARG	CONFLICT	GB 243868
A	100	SER	GLN	CONFLICT	GB 243868
A	104	LEU	VAL	CONFLICT	GB 243868

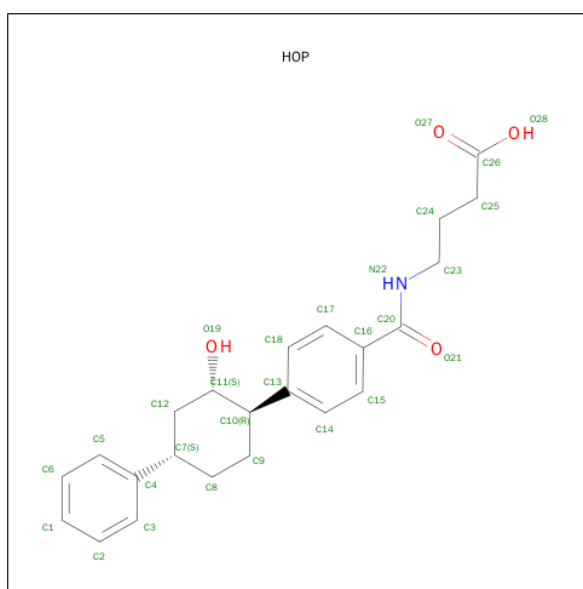
- Molecule 2 is a protein called OXY-COPE CATALYTIC ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1664	1057	271	330	6			
2	B	221	Total	C	N	O	S	0	0	0
			1664	1057	271	330	6			

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	2	Total	Cd	0	0
			2	2		
3	A	3	Total	Cd	0	0
			3	3		
3	L	3	Total	Cd	0	0
			3	3		

- Molecule 4 is (1S,2S,5S)2-(4-GLUTARIDYLBENZYL)-5-PHENYL-1-CYCLOHEXANOL (three-letter code: HOP) (formula: C₂₃H₂₇NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	1	Total	C	N	O	0	0
			28	23	1	4		
4	B	1	Total	C	N	O	0	0
			28	23	1	4		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	47	Total	O	0	0
			47	47		
5	B	37	Total	O	0	0
			37	37		
5	H	45	Total	O	0	0
			45	45		

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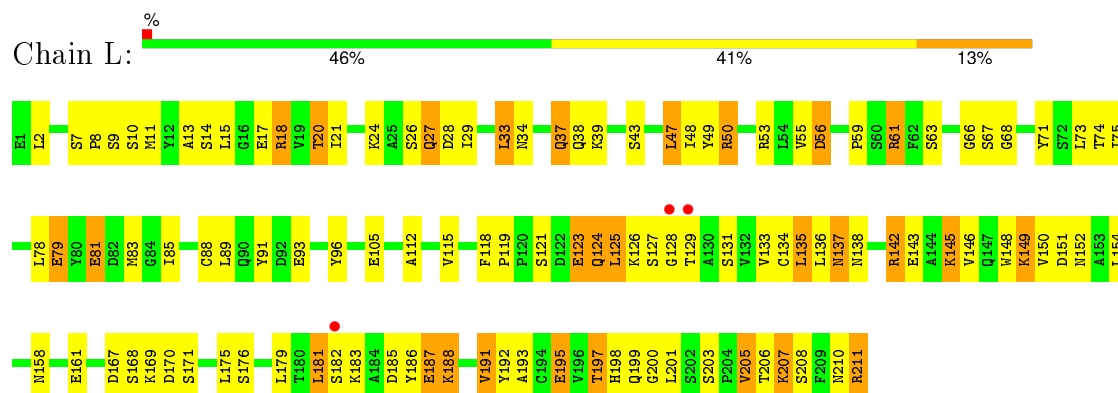
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	56	Total	O	0	0
			56	56		

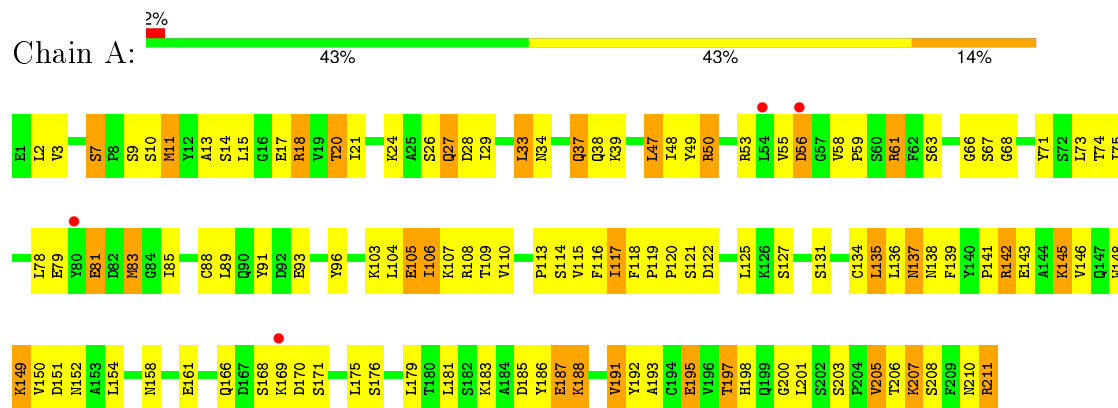
3 Residue-property plots

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

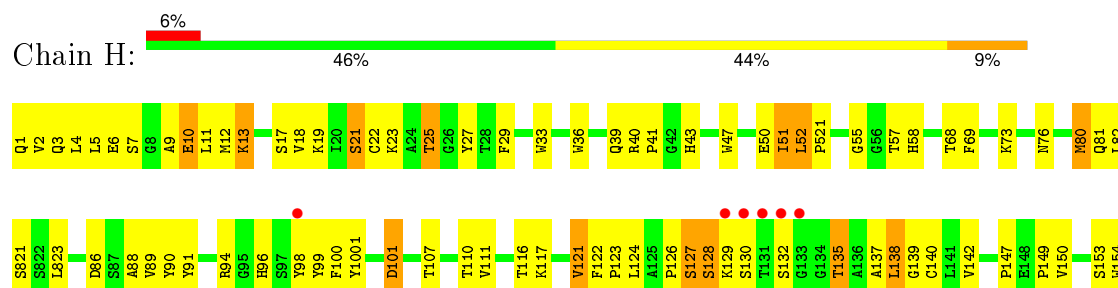
• Molecule 1: OXY-COPE CATALYTIC ANTIBODY



• Molecule 1: OXY-COPE CATALYTIC ANTIBODY

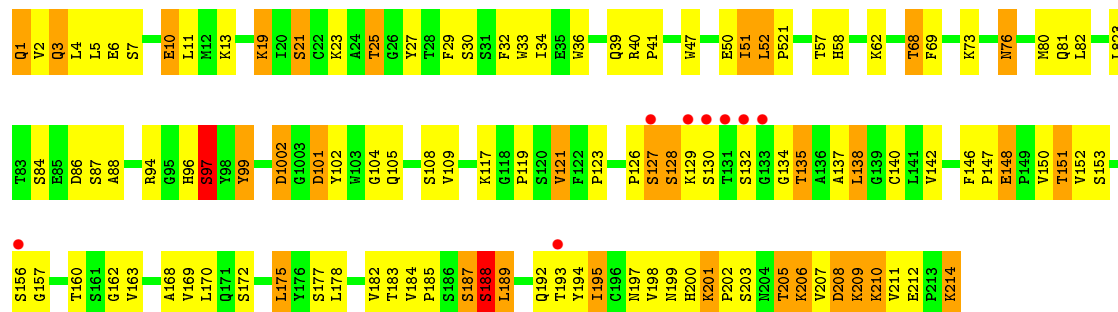


• Molecule 2: OXY-COPE CATALYTIC ANTIBODY





● Molecule 2: OXY-COPE CATALYTIC ANTIBODY



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.78 Å 81.50 Å 128.20 Å 90.00° 93.43° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 37.79 – 2.60	Depositor EDS
% Data completeness (in resolution range)	91.1 (20.00-2.60) 90.2 (37.79-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.17 (at 2.61 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.200 , 0.305 0.199 , 0.302	Depositor DCC
R_{free} test set	2434 reflections (9.92%)	DCC
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 72.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 24768 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6887	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: HOP, CD

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	0.39	0/1691	0.61	0/2291
1	L	0.45	1/1691 (0.1%)	0.64	0/2291
2	B	0.41	0/1710	0.66	2/2325 (0.1%)
2	H	0.41	0/1710	0.64	0/2325
All	All	0.41	1/6802 (0.0%)	0.64	2/9232 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	79	GLU	CB-CG	5.48	1.62	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	148	GLU	N-CA-C	5.73	126.48	111.00
2	B	97	SER	N-CA-C	-5.22	96.91	111.00

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	1655	0	1607	116	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1655	0	1607	96	0
2	B	1664	0	1610	103	0
2	H	1664	0	1610	83	0
3	A	3	0	0	0	0
3	H	2	0	0	0	0
3	L	3	0	0	0	0
4	B	28	0	27	4	0
4	H	28	0	27	4	0
5	A	47	0	0	11	0
5	B	37	0	0	12	0
5	H	45	0	0	5	0
5	L	56	0	0	6	0
All	All	6887	0	6488	384	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 29.

All (384) 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:L:38:GLN:HE22	2:H:39:GLN:HE22	1.04	1.00
1:L:137:ASN:HD22	1:L:138:ASN:N	1.60	0.99
1:A:58:VAL:HG21	5:A:215:HOH:O	1.63	0.99
2:B:33:TRP:HB3	5:B:1017:HOH:O	1.63	0.97
1:A:137:ASN:HD22	1:A:138:ASN:N	1.62	0.97
1:L:105:GLU:HG3	5:L:227:HOH:O	1.62	0.97
1:L:38:GLN:NE2	2:H:39:GLN:HE22	1.65	0.95
1:A:38:GLN:HE22	2:B:39:GLN:HE22	1.07	0.95
1:L:129:THR:HG22	1:L:182:SER:HA	1.50	0.93
1:A:15:LEU:HD11	1:A:108:ARG:HH22	1.30	0.93
2:B:50:GLU:HB2	5:B:1017:HOH:O	1.72	0.90
1:A:47:LEU:HA	5:A:215:HOH:O	1.71	0.89
1:A:38:GLN:NE2	2:B:39:GLN:HE22	1.72	0.86
1:A:55:VAL:HB	5:A:215:HOH:O	1.77	0.84
2:B:200:HIS:HB3	2:B:205:THR:HG23	1.60	0.84
1:L:38:GLN:HE22	2:H:39:GLN:NE2	1.74	0.84
1:A:210:ASN:O	1:A:211:ARG:HB2	1.77	0.83
2:B:135:THR:HG22	2:B:185:PRO:HA	1.60	0.83
1:L:210:ASN:O	1:L:211:ARG:HB2	1.77	0.83
2:B:156:SER:HA	2:B:197:ASN:HD21	1.44	0.82
2:B:156:SER:HA	2:B:197:ASN:ND2	1.95	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:135:THR:HG22	2:H:185:PRO:HA	1.60	0.81
2:B:156:SER:CA	2:B:197:ASN:HD21	1.97	0.77
1:A:38:GLN:HE22	2:B:39:GLN:NE2	1.82	0.76
1:L:137:ASN:HD22	1:L:138:ASN:H	1.30	0.76
1:A:137:ASN:HD22	1:A:138:ASN:H	1.32	0.75
2:B:1:GLN:HA	2:B:1:GLN:HE21	1.50	0.75
2:B:36:TRP:HE1	2:B:51:ILE:HD11	1.51	0.75
1:A:50:ARG:HG3	1:A:50:ARG:O	1.86	0.74
2:H:36:TRP:HE1	2:H:51:ILE:HD11	1.53	0.74
1:L:201:LEU:HD13	1:L:205:VAL:HG12	1.69	0.74
1:A:21:ILE:HB	5:A:247:HOH:O	1.88	0.73
1:L:50:ARG:HG3	1:L:50:ARG:O	1.87	0.72
1:A:201:LEU:HD13	1:A:205:VAL:HG12	1.70	0.72
1:A:20:THR:HB	1:A:74:THR:OG1	1.90	0.72
1:A:61:ARG:HD2	1:A:75:ILE:CG2	2.20	0.71
2:B:156:SER:N	2:B:197:ASN:HD21	1.88	0.71
2:H:201:LYS:O	2:H:203:SER:N	2.25	0.70
1:A:33:LEU:HD11	1:A:88:CYS:HB2	1.73	0.69
1:L:61:ARG:HD2	1:L:75:ILE:CG2	2.21	0.69
2:B:156:SER:H	2:B:197:ASN:HD21	1.40	0.69
2:B:195:ILE:HA	2:B:210:LYS:HA	1.75	0.68
1:L:20:THR:HB	1:L:74:THR:OG1	1.93	0.67
1:L:8:PRO:HD2	5:L:233:HOH:O	1.95	0.67
1:A:79:GLU:HG3	5:A:235:HOH:O	1.95	0.66
1:L:91:TYR:CD1	4:H:1006:HOP:H92	2.30	0.66
2:B:160:THR:O	2:B:163:VAL:HG23	1.96	0.66
1:L:161:GLU:OE2	1:L:175:LEU:HD11	1.96	0.66
1:A:28:ASP:OD2	5:A:258:HOH:O	2.13	0.66
1:L:145:LYS:HB3	1:L:197:THR:HG23	1.78	0.66
2:B:147:PRO:HD2	2:B:202:PRO:CB	2.25	0.65
1:A:158:ASN:ND2	1:A:179:LEU:HD11	2.11	0.65
1:A:91:TYR:CD1	4:B:1004:HOP:H92	2.32	0.65
1:L:158:ASN:ND2	1:L:179:LEU:HD11	2.11	0.64
2:B:198:VAL:C	2:B:199:ASN:HD22	2.02	0.64
1:A:161:GLU:OE2	1:A:175:LEU:HD11	1.97	0.64
1:A:145:LYS:HB3	1:A:197:THR:HG23	1.79	0.63
1:L:33:LEU:HD11	1:L:88:CYS:HB2	1.80	0.63
2:H:40:ARG:HG2	2:H:88:ALA:HB2	1.80	0.63
1:A:48:ILE:HD12	1:A:73:LEU:HD13	1.81	0.63
1:L:48:ILE:HD12	1:L:73:LEU:HD13	1.82	0.62
2:B:101:ASP:CG	4:B:1004:HOP:H3	2.20	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:ILE:O	2:B:195:ILE:HG13	2.00	0.62
1:A:15:LEU:HD12	1:A:108:ARG:HH12	1.64	0.62
2:B:32:PHE:CE1	2:B:97:SER:HB3	2.34	0.62
2:B:168:ALA:HA	2:B:178:LEU:HB3	1.80	0.62
1:L:201:LEU:HD13	1:L:205:VAL:CG1	2.29	0.62
2:H:195:ILE:HD12	2:H:197:ASN:HD21	1.65	0.61
1:L:136:LEU:HD13	1:L:175:LEU:HD22	1.82	0.61
2:B:51:ILE:HG12	2:B:69:PHE:CB	2.30	0.61
1:A:183:LYS:O	1:A:187:GLU:HG2	2.01	0.61
1:L:50:ARG:CG	1:L:50:ARG:HH11	2.12	0.61
1:L:81:GLU:OE2	5:L:268:HOH:O	2.15	0.61
1:A:50:ARG:HH11	1:A:50:ARG:CG	2.12	0.61
1:L:21:ILE:HD11	1:L:73:LEU:HD23	1.82	0.61
1:A:61:ARG:HG2	1:A:61:ARG:NH1	2.15	0.61
2:H:51:ILE:HG12	2:H:69:PHE:CB	2.31	0.60
1:L:14:SER:OG	1:L:17:GLU:HG3	2.02	0.60
2:H:51:ILE:HG23	2:H:57:THR:HG22	1.83	0.60
2:H:4:LEU:HD23	2:H:22:CYS:SG	2.42	0.60
2:B:40:ARG:HG2	2:B:88:ALA:HB2	1.83	0.60
1:L:183:LYS:O	1:L:187:GLU:HG2	2.02	0.60
2:H:168:ALA:HA	2:H:178:LEU:HB3	1.84	0.60
1:A:106:ILE:HG13	1:A:171:SER:CB	2.31	0.59
1:A:106:ILE:HG13	1:A:171:SER:HB3	1.84	0.59
1:L:61:ARG:NH1	1:L:61:ARG:HG2	2.17	0.59
1:L:123:GLU:O	1:L:126:LYS:HG2	2.03	0.59
1:L:55:VAL:HG12	1:L:56:ASP:N	2.18	0.59
1:A:91:TYR:CG	4:B:1004:HOP:H92	2.38	0.59
1:A:106:ILE:HG13	1:A:171:SER:OG	2.03	0.59
1:A:21:ILE:HD11	1:A:73:LEU:HD23	1.84	0.59
2:H:153:SER:O	2:H:196:CYS:HA	2.02	0.59
1:A:201:LEU:HD13	1:A:205:VAL:CG1	2.32	0.58
1:A:28:ASP:OD1	1:A:68:GLY:HA2	2.02	0.58
2:B:29:PHE:CE2	2:B:521:PRO:HB3	2.37	0.58
2:H:201:LYS:HG3	2:H:202:PRO:N	2.17	0.58
2:B:206:LYS:HG2	2:B:208:ASP:OD1	2.04	0.58
2:H:116:THR:HG23	5:H:1021:HOH:O	2.02	0.58
1:L:18:ARG:HH11	1:L:18:ARG:HB2	1.68	0.58
2:H:123:PRO:HB2	2:H:211:VAL:HG13	1.83	0.58
2:H:9:ALA:C	2:H:10:GLU:HG2	2.22	0.58
1:L:125:LEU:CD2	1:L:183:LYS:HG3	2.33	0.58
1:A:18:ARG:HB2	1:A:18:ARG:HH11	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:TRP:NE1	2:H:51:ILE:HD11	2.19	0.57
1:L:15:LEU:HA	1:L:78:LEU:O	2.03	0.57
2:H:29:PHE:O	2:H:521:PRO:HG2	2.03	0.57
2:B:193:THR:HG22	2:B:194:TYR:H	1.69	0.57
2:B:51:ILE:HG23	2:B:57:THR:HG22	1.85	0.57
2:H:13:LYS:HG2	5:H:1051:HOH:O	2.05	0.57
1:A:55:VAL:HG12	1:A:56:ASP:N	2.20	0.57
1:A:15:LEU:HD21	1:A:106:ILE:HG21	1.86	0.57
2:B:36:TRP:NE1	2:B:51:ILE:HD11	2.19	0.57
1:A:14:SER:OG	1:A:17:GLU:HG3	2.05	0.57
1:A:106:ILE:HG12	1:A:166:GLN:NE2	2.20	0.56
1:A:15:LEU:HD11	1:A:108:ARG:NH2	2.11	0.56
1:L:91:TYR:CG	4:H:1006:HOP:H92	2.41	0.56
1:L:28:ASP:OD1	1:L:68:GLY:HA2	2.05	0.56
1:A:136:LEU:HD13	1:A:175:LEU:HD22	1.86	0.56
1:L:66:GLY:HA3	1:L:71:TYR:HA	1.88	0.56
2:H:127:SER:O	2:H:129:LYS:N	2.38	0.56
1:A:61:ARG:HG2	1:A:61:ARG:HH11	1.70	0.56
1:A:149:LYS:HB2	1:A:193:ALA:HB3	1.88	0.56
2:B:127:SER:O	2:B:129:LYS:N	2.39	0.56
1:A:50:ARG:CG	1:A:50:ARG:O	2.53	0.55
1:A:15:LEU:CD1	1:A:108:ARG:HH22	2.11	0.55
1:L:38:GLN:CD	2:H:39:GLN:HE22	2.09	0.55
1:A:210:ASN:O	1:A:211:ARG:HD2	2.06	0.55
2:H:184:VAL:HG11	2:H:194:TYR:CE2	2.42	0.55
2:B:11:LEU:HB2	2:B:147:PRO:HG3	1.88	0.55
1:A:66:GLY:HA3	1:A:71:TYR:HA	1.89	0.55
2:H:94:ARG:O	2:H:101:ASP:HB3	2.07	0.55
1:L:50:ARG:CG	1:L:50:ARG:O	2.55	0.54
1:L:151:ASP:O	1:L:152:ASN:HB2	2.07	0.54
1:L:149:LYS:HB2	1:L:193:ALA:HB3	1.89	0.54
2:H:82:LEU:HB2	2:H:823:LEU:HD21	1.89	0.54
1:A:15:LEU:HA	1:A:78:LEU:O	2.07	0.54
1:L:61:ARG:HG2	1:L:61:ARG:HH11	1.72	0.54
2:B:823:LEU:HA	2:B:86:ASP:OD2	2.08	0.54
1:L:198:HIS:CD2	1:L:200:GLY:H	2.26	0.53
2:H:33:TRP:CE2	2:H:52:LEU:HG	2.43	0.53
1:L:121:SER:OG	1:L:124:GLN:HB3	2.07	0.53
2:B:36:TRP:HE1	2:B:51:ILE:CD1	2.21	0.53
2:B:137:ALA:HB2	2:B:183:THR:HG22	1.91	0.53
2:B:33:TRP:CE2	2:B:52:LEU:HG	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:91:TYR:HA	1:L:96:TYR:CD1	2.44	0.53
1:L:185:ASP:HA	1:L:188:LYS:HD3	1.90	0.53
1:L:210:ASN:O	1:L:211:ARG:HD2	2.09	0.53
2:H:36:TRP:HE1	2:H:51:ILE:CD1	2.21	0.53
1:L:136:LEU:HD13	1:L:175:LEU:CD2	2.38	0.53
1:A:91:TYR:HA	1:A:96:TYR:CD1	2.43	0.53
1:L:79:GLU:HB3	1:L:81:GLU:OE2	2.10	0.52
2:H:194:TYR:N	2:H:194:TYR:CD1	2.77	0.52
1:A:185:ASP:HA	1:A:188:LYS:HD3	1.90	0.52
2:H:11:LEU:HD12	2:H:110:THR:O	2.09	0.52
2:B:147:PRO:HD2	2:B:202:PRO:HB3	1.91	0.52
1:L:27:GLN:O	1:L:29:ILE:HG23	2.09	0.52
2:B:6:GLU:O	2:B:7:SER:HB2	2.08	0.52
1:A:198:HIS:CD2	1:A:200:GLY:H	2.27	0.52
1:L:151:ASP:OD1	1:L:191:VAL:HG23	2.10	0.52
2:B:6:GLU:OE2	2:B:104:GLY:HA3	2.09	0.52
1:L:112:ALA:HB1	1:L:201:LEU:CD2	2.40	0.52
1:L:50:ARG:CG	1:L:50:ARG:NH1	2.71	0.52
1:A:79:GLU:HG2	5:A:260:HOH:O	2.09	0.52
1:A:210:ASN:O	1:A:211:ARG:CB	2.54	0.52
2:H:11:LEU:HD22	2:H:147:PRO:HG3	1.91	0.52
2:H:200:HIS:CE1	2:H:202:PRO:HB2	2.45	0.52
1:L:193:ALA:HA	1:L:208:SER:HB3	1.91	0.51
2:H:126:PRO:HB3	2:H:138:LEU:HB3	1.92	0.51
2:B:150:VAL:HG22	2:B:151:THR:N	2.24	0.51
2:H:137:ALA:HB2	2:H:183:THR:HG22	1.92	0.51
1:A:106:ILE:HG12	1:A:166:GLN:HE22	1.74	0.51
2:B:214:LYS:HG3	5:B:1030:HOH:O	2.10	0.51
2:H:139:GLY:HA2	2:H:154:TRP:CH2	2.46	0.51
2:H:96:HIS:HB2	2:H:1001:TYR:HB2	1.92	0.51
2:B:1:GLN:HA	2:B:1:GLN:NE2	2.22	0.51
1:A:146:VAL:HA	1:A:195:GLU:O	2.10	0.51
1:A:108:ARG:HG2	1:A:109:THR:H	1.75	0.51
2:B:200:HIS:HB3	2:B:205:THR:CG2	2.34	0.51
2:H:139:GLY:C	2:H:154:TRP:HH2	2.14	0.51
1:A:108:ARG:HE	1:A:171:SER:HB2	1.76	0.51
1:A:50:ARG:CG	1:A:50:ARG:NH1	2.72	0.51
1:A:192:TYR:O	1:A:208:SER:HA	2.11	0.51
1:A:83:MET:HE3	1:A:105:GLU:HA	1.92	0.51
2:B:82:LEU:HB2	2:B:823:LEU:HD21	1.92	0.51
1:A:27:GLN:O	1:A:29:ILE:HG23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:47:TRP:HZ2	2:B:50:GLU:HG2	1.75	0.51
2:H:147:PRO:HD2	2:H:202:PRO:HB3	1.93	0.51
2:B:1:GLN:HE21	2:B:1:GLN:CA	2.16	0.51
1:A:151:ASP:OD1	1:A:191:VAL:HG23	2.11	0.51
1:A:151:ASP:O	1:A:152:ASN:HB2	2.11	0.51
2:H:98:TYR:CE1	2:H:99:TYR:HD1	2.29	0.51
2:H:147:PRO:HD2	2:H:202:PRO:CB	2.41	0.50
1:A:15:LEU:CD1	1:A:108:ARG:HH12	2.24	0.50
2:B:203:SER:OG	2:B:205:THR:HG22	2.11	0.50
1:A:61:ARG:CG	1:A:61:ARG:HH11	2.25	0.50
2:H:210:LYS:HE3	2:H:212:GLU:OE2	2.12	0.50
1:A:136:LEU:HD13	1:A:175:LEU:CD2	2.41	0.50
1:L:133:VAL:HG23	5:L:228:HOH:O	2.11	0.50
2:H:101:ASP:CG	4:H:1006:HOP:H3	2.32	0.50
1:A:193:ALA:HA	1:A:208:SER:HB3	1.93	0.50
1:L:112:ALA:HB1	1:L:201:LEU:HD21	1.94	0.49
2:B:126:PRO:HB3	2:B:138:LEU:HB3	1.93	0.49
1:L:146:VAL:HA	1:L:195:GLU:O	2.12	0.49
1:A:170:ASP:O	1:A:171:SER:HB2	2.12	0.49
2:B:200:HIS:HD2	2:B:203:SER:OG	1.95	0.49
1:A:14:SER:HB2	5:A:221:HOH:O	2.12	0.49
1:L:192:TYR:O	1:L:208:SER:HA	2.13	0.49
2:B:178:LEU:C	2:B:178:LEU:HD12	2.32	0.49
1:A:135:LEU:HD22	1:A:136:LEU:N	2.28	0.49
2:H:126:PRO:HD3	2:H:138:LEU:HB2	1.93	0.49
1:A:121:SER:HB2	2:B:123:PRO:HD2	1.94	0.49
2:B:169:VAL:HG12	2:B:177:SER:O	2.13	0.49
2:H:29:PHE:CE2	2:H:521:PRO:HB3	2.48	0.49
2:B:10:GLU:HG3	5:B:1032:HOH:O	2.12	0.49
1:A:125:LEU:O	1:A:127:SER:O	2.31	0.49
1:A:108:ARG:NE	1:A:171:SER:HB2	2.28	0.49
2:H:47:TRP:HZ2	2:H:50:GLU:HG2	1.78	0.49
1:A:134:CYS:HB2	1:A:148:TRP:CZ2	2.48	0.49
2:H:154:TRP:O	2:H:155:ASN:HB2	2.13	0.48
1:A:79:GLU:HB3	1:A:81:GLU:OE2	2.13	0.48
1:L:170:ASP:O	1:L:171:SER:HB2	2.13	0.48
2:H:178:LEU:C	2:H:178:LEU:HD12	2.33	0.48
2:B:153:SER:OG	2:B:197:ASN:HB2	2.14	0.48
2:B:3:GLN:HB2	2:B:25:THR:HG22	1.96	0.48
1:L:135:LEU:HD22	1:L:136:LEU:N	2.28	0.48
1:A:61:ARG:HD2	1:A:75:ILE:HG23	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:61:ARG:CG	1:L:61:ARG:HH11	2.27	0.48
2:B:6:GLU:HA	2:B:21:SER:O	2.14	0.48
2:B:126:PRO:HD3	2:B:138:LEU:HB2	1.95	0.48
2:B:193:THR:HG22	2:B:194:TYR:N	2.29	0.47
1:L:61:ARG:HD2	1:L:75:ILE:HG23	1.96	0.47
2:H:43:HIS:HB2	5:H:1019:HOH:O	2.13	0.47
2:B:1:GLN:CA	2:B:1:GLN:NE2	2.78	0.47
1:A:113:PRO:HB3	1:A:139:PHE:HB3	1.96	0.47
2:B:76:ASN:ND2	5:B:1027:HOH:O	2.40	0.47
1:A:116:PHE:HA	2:B:129:LYS:NZ	2.30	0.47
1:L:134:CYS:HB2	1:L:148:TRP:CZ2	2.49	0.47
2:B:84:SER:HA	5:B:1028:HOH:O	2.13	0.47
1:L:50:ARG:HG2	1:L:53:ARG:CG	2.45	0.47
2:B:27:TYR:CE2	2:B:94:ARG:HD2	2.50	0.47
1:L:55:VAL:CG1	1:L:56:ASP:N	2.77	0.47
1:A:142:ARG:HG3	1:A:143:GLU:N	2.29	0.47
2:B:201:LYS:HB3	2:B:202:PRO:HD3	1.97	0.47
2:H:187:SER:O	2:H:189:LEU:N	2.48	0.47
2:B:30:SER:HA	2:B:521:PRO:HB2	1.95	0.47
1:L:37:GLN:HB3	1:L:47:LEU:HD22	1.96	0.47
1:L:124:GLN:HB2	2:H:122:PHE:CE2	2.50	0.47
2:H:51:ILE:CD1	2:H:69:PHE:HB3	2.45	0.47
2:H:150:VAL:HG12	5:H:1014:HOH:O	2.14	0.47
2:H:52:LEU:HD23	2:H:52:LEU:HA	1.82	0.46
1:A:154:LEU:O	1:A:154:LEU:HD23	2.14	0.46
1:L:142:ARG:HG3	1:L:143:GLU:N	2.30	0.46
1:L:129:THR:HB	1:L:181:LEU:O	2.15	0.46
1:L:61:ARG:O	1:L:75:ILE:HA	2.15	0.46
1:L:154:LEU:HD23	1:L:154:LEU:O	2.14	0.46
2:H:121:VAL:CG1	2:H:207:VAL:HG11	2.45	0.46
2:H:823:LEU:HA	2:H:86:ASP:OD2	2.15	0.46
1:L:137:ASN:ND2	1:L:138:ASN:N	2.44	0.46
1:A:55:VAL:CG1	1:A:56:ASP:N	2.78	0.46
2:B:184:VAL:HG11	2:B:194:TYR:CZ	2.51	0.46
1:L:210:ASN:O	1:L:211:ARG:CB	2.55	0.46
1:A:119:PRO:O	1:A:120:PRO:C	2.54	0.46
2:B:187:SER:O	2:B:189:LEU:N	2.48	0.46
2:H:7:SER:HB3	2:H:21:SER:OG	2.15	0.46
1:A:61:ARG:O	1:A:75:ILE:HA	2.16	0.46
1:L:175:LEU:HD23	1:L:175:LEU:C	2.36	0.46
2:H:11:LEU:HD23	2:H:116:THR:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:MET:HB3	1:A:104:LEU:CD1	2.45	0.45
2:H:206:LYS:HE3	2:H:206:LYS:HB2	1.61	0.45
1:A:37:GLN:HB3	1:A:47:LEU:HD22	1.97	0.45
2:B:121:VAL:CG1	2:B:207:VAL:HG11	2.46	0.45
2:H:27:TYR:CE2	2:H:94:ARG:HD2	2.51	0.45
1:A:107:LYS:HE3	5:A:240:HOH:O	2.14	0.45
1:L:48:ILE:HD12	1:L:73:LEU:CD1	2.46	0.45
1:A:175:LEU:HD23	1:A:175:LEU:C	2.36	0.45
2:B:68:THR:HG22	5:B:1024:HOH:O	2.15	0.45
1:A:26:SER:O	1:A:27:GLN:HB3	2.15	0.45
2:H:199:ASN:OD1	2:H:206:LYS:HG3	2.16	0.45
1:L:34:ASN:OD1	1:L:49:TYR:HA	2.15	0.45
1:A:38:GLN:CD	2:B:39:GLN:HE22	2.19	0.45
2:B:195:ILE:HD11	2:B:197:ASN:OD1	2.16	0.45
2:B:51:ILE:CD1	2:B:69:PHE:HB3	2.47	0.45
1:L:13:ALA:HB3	1:L:78:LEU:HD22	1.97	0.45
2:B:123:PRO:HB2	2:B:211:VAL:HG13	1.99	0.45
1:A:50:ARG:HG2	1:A:53:ARG:CG	2.46	0.45
2:B:146:PHE:HB2	2:B:175:LEU:HD12	1.98	0.45
2:B:47:TRP:CZ2	2:B:50:GLU:HG2	2.52	0.44
1:L:150:VAL:HG22	1:L:192:TYR:CD1	2.53	0.44
2:H:169:VAL:HG12	2:H:177:SER:O	2.16	0.44
2:H:13:LYS:HG2	2:H:13:LYS:H	1.42	0.44
1:A:50:ARG:HH11	1:A:50:ARG:HG2	1.81	0.44
1:L:50:ARG:HG3	1:L:50:ARG:NH1	2.31	0.44
1:A:117:ILE:HG13	1:A:118:PHE:N	2.33	0.44
1:L:50:ARG:HG2	1:L:53:ARG:HG3	2.00	0.44
1:L:183:LYS:O	1:L:186:TYR:HB3	2.18	0.44
2:H:142:VAL:CG2	2:H:198:VAL:HG21	2.48	0.44
1:A:137:ASN:ND2	1:A:138:ASN:N	2.46	0.44
2:H:2:VAL:HA	2:H:25:THR:O	2.18	0.44
2:B:2:VAL:HA	2:B:25:THR:O	2.17	0.44
2:H:101:ASP:OD1	4:H:1006:HOP:H7	2.17	0.44
1:A:113:PRO:HB3	1:A:139:PHE:CG	2.53	0.44
2:B:34:ILE:N	5:B:1017:HOH:O	2.51	0.43
2:B:152:VAL:HA	2:B:197:ASN:O	2.17	0.43
1:A:150:VAL:HG22	1:A:192:TYR:CD1	2.53	0.43
1:A:193:ALA:CB	1:A:208:SER:HB3	2.49	0.43
2:H:51:ILE:HG12	2:H:69:PHE:CG	2.53	0.43
1:L:27:GLN:O	1:L:28:ASP:C	2.56	0.43
2:B:19:LYS:HB2	2:B:19:LYS:HE3	1.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:185:PRO:C	2:B:187:SER:N	2.71	0.43
2:H:185:PRO:C	2:H:187:SER:N	2.71	0.43
1:L:26:SER:O	1:L:27:GLN:HB3	2.18	0.43
2:B:187:SER:OG	2:B:188:SER:N	2.52	0.43
2:B:2:VAL:HB	2:B:102:TYR:CD2	2.53	0.43
1:A:110:VAL:HG22	1:A:141:PRO:HD3	2.00	0.43
2:B:13:LYS:HG3	5:B:1006:HOH:O	2.17	0.43
2:H:123:PRO:CB	2:H:211:VAL:HG13	2.49	0.43
2:B:200:HIS:CD2	2:B:203:SER:OG	2.72	0.43
2:B:209:LYS:HD2	2:B:210:LYS:N	2.33	0.43
2:B:51:ILE:HG12	2:B:69:PHE:CG	2.54	0.43
2:H:201:LYS:C	2:H:203:SER:H	2.22	0.43
2:H:124:LEU:HB2	2:H:139:GLY:C	2.39	0.43
1:A:206:THR:O	1:A:207:LYS:HD2	2.19	0.43
2:B:207:VAL:HG12	2:B:207:VAL:O	2.19	0.42
1:A:48:ILE:HD12	1:A:73:LEU:CD1	2.46	0.42
1:A:61:ARG:HD2	1:A:75:ILE:HG22	1.98	0.42
1:L:112:ALA:HB2	1:L:200:GLY:O	2.18	0.42
1:L:115:VAL:HA	1:L:135:LEU:O	2.20	0.42
1:L:193:ALA:CA	1:L:208:SER:HB3	2.49	0.42
2:B:184:VAL:HG11	2:B:194:TYR:CE2	2.54	0.42
2:B:62:LYS:HG3	5:B:1018:HOH:O	2.19	0.42
1:A:34:ASN:OD1	1:A:49:TYR:HA	2.19	0.42
2:B:147:PRO:O	2:B:200:HIS:HE1	2.01	0.42
1:L:61:ARG:HD2	1:L:75:ILE:HG22	2.00	0.42
1:A:50:ARG:HG3	1:A:50:ARG:NH1	2.33	0.42
1:L:187:GLU:HG2	1:L:187:GLU:H	1.50	0.42
1:A:7:SER:HB3	5:A:250:HOH:O	2.19	0.42
2:H:12:MET:O	2:H:111:VAL:HA	2.19	0.42
1:A:49:TYR:CE1	2:B:1002:ASP:O	2.73	0.42
2:H:3:GLN:HG2	5:H:1010:HOH:O	2.19	0.42
1:A:13:ALA:HB3	1:A:78:LEU:HD22	2.00	0.42
1:A:193:ALA:CA	1:A:208:SER:HB3	2.49	0.42
1:A:113:PRO:HB3	1:A:139:PHE:CB	2.50	0.42
2:B:119:PRO:HB2	2:B:142:VAL:HG13	2.02	0.42
1:L:50:ARG:HH11	1:L:50:ARG:HG2	1.83	0.42
2:B:193:THR:HB	5:B:1007:HOH:O	2.20	0.42
2:H:100:PHE:CG	2:H:1001:TYR:N	2.89	0.41
1:L:206:THR:O	1:L:207:LYS:HD2	2.20	0.41
2:B:207:VAL:O	2:B:209:LYS:N	2.53	0.41
2:H:69:PHE:CD1	2:H:80:MET:HG3	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:154:TRP:HA	2:H:195:ILE:O	2.21	0.41
1:A:113:PRO:CA	1:A:139:PHE:HB3	2.49	0.41
2:H:89:VAL:HA	2:H:107:THR:O	2.20	0.41
1:L:43:SER:HB2	2:H:91:TYR:CE1	2.55	0.41
1:L:124:GLN:NE2	1:L:129:THR:O	2.54	0.41
1:A:183:LYS:O	1:A:186:TYR:HB3	2.20	0.41
1:A:113:PRO:HD3	1:A:198:HIS:CD2	2.55	0.41
1:A:3:VAL:HA	5:A:242:HOH:O	2.19	0.41
2:B:87:SER:HA	2:B:109:VAL:O	2.20	0.41
2:H:187:SER:OG	2:H:188:SER:N	2.53	0.41
2:B:163:VAL:HG22	2:B:182:VAL:HG23	2.02	0.41
1:A:193:ALA:HB2	1:A:208:SER:HB3	2.02	0.41
2:B:1:GLN:HG3	2:B:2:VAL:O	2.20	0.41
2:H:52:LEU:O	2:H:55:GLY:N	2.50	0.41
2:H:47:TRP:CZ2	2:H:50:GLU:HG2	2.55	0.41
1:A:50:ARG:HG2	1:A:53:ARG:HG3	2.03	0.41
2:B:1002:ASP:N	5:B:1019:HOH:O	2.53	0.41
1:L:118:PHE:HA	1:L:119:PRO:HD3	1.81	0.41
1:L:150:VAL:O	1:L:151:ASP:HB2	2.21	0.41
2:B:99:TYR:CD1	2:B:99:TYR:C	2.94	0.41
2:B:96:HIS:ND1	4:B:1004:HOP:O19	2.41	0.41
1:L:193:ALA:CB	1:L:208:SER:HB3	2.50	0.41
2:H:50:GLU:HG3	2:H:58:HIS:HB2	2.03	0.41
2:B:162:GLY:O	2:B:182:VAL:HA	2.21	0.41
1:L:37:GLN:HG2	5:L:223:HOH:O	2.21	0.41
2:H:164:HIS:HB2	2:H:181:VAL:CG2	2.51	0.41
2:B:195:ILE:HG22	2:B:210:LYS:HB2	2.03	0.40
1:L:167:ASP:O	1:L:171:SER:N	2.49	0.40
1:A:121:SER:CB	2:B:123:PRO:HD2	2.52	0.40
1:L:61:ARG:HB3	5:L:218:HOH:O	2.21	0.40
1:A:115:VAL:HA	1:A:135:LEU:O	2.21	0.40
2:H:18:VAL:HB	2:H:823:LEU:HD11	2.04	0.40
2:B:50:GLU:HG3	2:B:58:HIS:HB2	2.02	0.40
2:H:6:GLU:OE1	2:H:90:TYR:O	2.39	0.40
1:A:83:MET:HG3	1:A:106:ILE:CD1	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	209/211 (99%)	191 (91%)	17 (8%)	1 (0%)	34	60
1	L	209/211 (99%)	191 (91%)	15 (7%)	3 (1%)	14	28
2	B	219/221 (99%)	187 (85%)	23 (10%)	9 (4%)	3	4
2	H	219/221 (99%)	190 (87%)	21 (10%)	8 (4%)	4	5
All	All	856/864 (99%)	759 (89%)	76 (9%)	21 (2%)	7	12

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	188	SER
2	H	202	PRO
2	H	213	PRO
1	L	128	GLY
2	H	128	SER
2	B	128	SER
2	B	188	SER
2	H	187	SER
2	B	187	SER
2	H	132	SER
2	B	132	SER
2	B	208	ASP
2	H	41	PRO
2	B	41	PRO
2	B	157	GLY
1	L	199	GLN
2	B	99	TYR
2	H	149	PRO
1	A	59	PRO
1	L	59	PRO
2	B	134	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	190/190 (100%)	142 (75%)	48 (25%)	1	1
1	L	190/190 (100%)	144 (76%)	46 (24%)	1	1
2	B	185/185 (100%)	140 (76%)	45 (24%)	1	1
2	H	185/185 (100%)	145 (78%)	40 (22%)	1	2
All	All	750/750 (100%)	571 (76%)	179 (24%)	1	1

All (179) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	LEU
1	L	7	SER
1	L	9	SER
1	L	10	SER
1	L	11	MET
1	L	18	ARG
1	L	20	THR
1	L	24	LYS
1	L	27	GLN
1	L	33	LEU
1	L	37	GLN
1	L	39	LYS
1	L	47	LEU
1	L	50	ARG
1	L	56	ASP
1	L	61	ARG
1	L	63	SER
1	L	67	SER
1	L	81	GLU
1	L	83	MET
1	L	85	ILE
1	L	89	LEU
1	L	93	GLU
1	L	123	GLU

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Mol	Chain	Res	Type
1	L	124	GLN
1	L	125	LEU
1	L	127	SER
1	L	131	SER
1	L	135	LEU
1	L	137	ASN
1	L	142	ARG
1	L	145	LYS
1	L	149	LYS
1	L	168	SER
1	L	169	LYS
1	L	176	SER
1	L	181	LEU
1	L	187	GLU
1	L	188	LYS
1	L	191	VAL
1	L	195	GLU
1	L	197	THR
1	L	203	SER
1	L	205	VAL
1	L	207	LYS
1	L	211	ARG
2	H	1	GLN
2	H	5	LEU
2	H	10	GLU
2	H	13	LYS
2	H	17	SER
2	H	19	LYS
2	H	21	SER
2	H	23	LYS
2	H	25	THR
2	H	51	ILE
2	H	52	LEU
2	H	68	THR
2	H	73	LYS
2	H	76	ASN
2	H	80	MET
2	H	81	GLN
2	H	821	SER
2	H	101	ASP
2	H	117	LYS
2	H	121	VAL

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Mol	Chain	Res	Type
2	H	127	SER
2	H	128	SER
2	H	130	SER
2	H	135	THR
2	H	138	LEU
2	H	140	CYS
2	H	161	SER
2	H	167	PRO
2	H	170	LEU
2	H	175	LEU
2	H	188	SER
2	H	189	LEU
2	H	192	GLN
2	H	194	TYR
2	H	195	ILE
2	H	201	LYS
2	H	205	THR
2	H	209	LYS
2	H	211	VAL
2	H	214	LYS
1	A	2	LEU
1	A	7	SER
1	A	9	SER
1	A	10	SER
1	A	11	MET
1	A	18	ARG
1	A	20	THR
1	A	24	LYS
1	A	27	GLN
1	A	33	LEU
1	A	37	GLN
1	A	39	LYS
1	A	47	LEU
1	A	50	ARG
1	A	56	ASP
1	A	61	ARG
1	A	63	SER
1	A	67	SER
1	A	81	GLU
1	A	83	MET
1	A	85	ILE
1	A	89	LEU

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Mol	Chain	Res	Type
1	A	93	GLU
1	A	103	LYS
1	A	105	GLU
1	A	106	ILE
1	A	114	SER
1	A	117	ILE
1	A	122	ASP
1	A	131	SER
1	A	135	LEU
1	A	137	ASN
1	A	142	ARG
1	A	145	LYS
1	A	149	LYS
1	A	168	SER
1	A	169	LYS
1	A	176	SER
1	A	181	LEU
1	A	187	GLU
1	A	188	LYS
1	A	191	VAL
1	A	195	GLU
1	A	197	THR
1	A	203	SER
1	A	205	VAL
1	A	207	LYS
1	A	211	ARG
2	B	1	GLN
2	B	3	GLN
2	B	4	LEU
2	B	5	LEU
2	B	10	GLU
2	B	19	LYS
2	B	21	SER
2	B	23	LYS
2	B	25	THR
2	B	51	ILE
2	B	52	LEU
2	B	68	THR
2	B	73	LYS
2	B	76	ASN
2	B	80	MET
2	B	81	GLN

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Mol	Chain	Res	Type
2	B	97	SER
2	B	1002	ASP
2	B	101	ASP
2	B	105	GLN
2	B	108	SER
2	B	117	LYS
2	B	121	VAL
2	B	127	SER
2	B	128	SER
2	B	130	SER
2	B	135	THR
2	B	138	LEU
2	B	140	CYS
2	B	148	GLU
2	B	151	THR
2	B	170	LEU
2	B	172	SER
2	B	175	LEU
2	B	188	SER
2	B	189	LEU
2	B	192	GLN
2	B	195	ILE
2	B	201	LYS
2	B	205	THR
2	B	206	LYS
2	B	209	LYS
2	B	210	LYS
2	B	212	GLU
2	B	214	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	GLN
1	L	69	GLN
1	L	124	GLN
1	L	137	ASN
1	L	198	HIS
2	H	39	GLN
2	H	43	HIS
2	H	58	HIS
1	A	37	GLN

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Mol	Chain	Res	Type
1	A	69	GLN
1	A	137	ASN
1	A	198	HIS
2	B	1	GLN
2	B	3	GLN
2	B	39	GLN
2	B	43	HIS
2	B	58	HIS
2	B	197	ASN
2	B	199	ASN
2	B	200	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 8 are monoatomic - leaving 2 for Mogul analysis.

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
4	HOP	B	1004	-	27,30,30	2.09	10 (37%)	35,40,40	0.62	1 (2%)
4	HOP	H	1006	-	27,30,30	2.36	15 (55%)	35,40,40	0.83	1 (2%)

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
4	HOP	B	1004	-	-	0/17/32/32	0/3/3/3
4	HOP	H	1006	-	-	0/17/32/32	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1006	HOP	C20-N22	2.04	1.38	1.33
4	H	1006	HOP	C3-C4	2.16	1.42	1.39
4	H	1006	HOP	C16-C20	2.23	1.54	1.50
4	H	1006	HOP	C17-C16	2.28	1.43	1.39
4	H	1006	HOP	C4-C7	2.34	1.56	1.52
4	H	1006	HOP	C18-C13	2.48	1.43	1.39
4	B	1004	HOP	C14-C13	2.48	1.43	1.39
4	B	1004	HOP	C3-C4	2.51	1.43	1.39
4	B	1004	HOP	C9-C10	2.58	1.57	1.53
4	H	1006	HOP	C12-C11	2.65	1.57	1.52
4	B	1004	HOP	C12-C7	2.68	1.56	1.53
4	B	1004	HOP	C17-C16	2.74	1.44	1.39
4	B	1004	HOP	C15-C16	2.75	1.44	1.39
4	B	1004	HOP	C18-C13	2.75	1.43	1.39
4	H	1006	HOP	C9-C10	3.00	1.57	1.53
4	H	1006	HOP	C10-C11	3.06	1.57	1.53
4	H	1006	HOP	C14-C13	3.08	1.44	1.39
4	H	1006	HOP	C15-C14	3.11	1.44	1.38
4	B	1004	HOP	C5-C4	3.27	1.44	1.39
4	H	1006	HOP	C5-C4	3.41	1.44	1.39
4	B	1004	HOP	C10-C11	3.44	1.57	1.53
4	H	1006	HOP	C12-C7	3.47	1.57	1.53
4	H	1006	HOP	C15-C16	3.62	1.45	1.39
4	H	1006	HOP	C13-C10	3.81	1.57	1.51
4	B	1004	HOP	C13-C10	3.87	1.57	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1006	HOP	C18-C13-C10	-2.24	116.59	121.08
4	B	1004	HOP	C9-C10-C13	-2.07	109.24	112.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1004	HOP	4	0
4	H	1006	HOP	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	211/211 (100%)	-0.01	4 (1%) 70 64	20, 42, 67, 87	0
1	L	211/211 (100%)	-0.25	3 (1%) 78 74	8, 27, 57, 71	0
2	B	221/221 (100%)	0.00	8 (3%) 46 38	9, 33, 82, 99	0
2	H	221/221 (100%)	0.05	14 (6%) 23 17	9, 31, 79, 100	0
All	All	864/864 (100%)	-0.05	29 (3%) 49 41	8, 33, 72, 100	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	131	THR	7.1
2	H	130	SER	6.4
2	B	127	SER	4.6
2	H	133	GLY	4.5
2	H	161	SER	4.4
2	H	132	SER	4.2
2	H	191	THR	4.1
2	B	133	GLY	3.8
2	H	193	THR	3.5
2	H	159	LEU	3.4
2	H	187	SER	3.3
1	A	169	LYS	3.1
1	L	129	THR	3.1
2	H	190	GLY	3.0
1	A	54	LEU	2.9
2	B	131	THR	2.9
2	H	172	SER	2.6
2	B	193	THR	2.6
2	B	130	SER	2.6
2	B	132	SER	2.5
1	L	128	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	129	LYS	2.5
1	A	56	ASP	2.4
1	L	182	SER	2.3
2	B	156	SER	2.2
2	H	188	SER	2.2
1	A	80	TYR	2.1
2	H	98	TYR	2.1
2	H	129	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	LLDF	B-factors(Å ²)	Q<0.9
4	HOP	H	1006	28/28	0.90	0.21	4.06	12,30,63,66	0
4	HOP	B	1004	28/28	0.89	0.22	1.35	26,50,71,82	0
3	CD	L	214	1/1	1.00	0.13	-0.80	27,27,27,27	0
3	CD	L	212	1/1	0.98	0.14	-1.03	39,39,39,39	1
3	CD	A	212	1/1	0.96	0.12	-1.16	47,47,47,47	1
3	CD	A	213	1/1	0.96	0.09	-1.39	55,55,55,55	1
3	CD	A	214	1/1	1.00	0.07	-3.40	46,46,46,46	0
3	CD	L	213	1/1	0.98	0.06	-3.71	57,57,57,57	0
3	CD	H	1004	1/1	0.98	0.05	-	50,50,50,50	1
3	CD	H	1005	1/1	0.95	0.07	-	85,85,85,85	0

6.5 Other polymers

There are no such residues in this entry.