



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:53 pm GMT

PDB ID : 2F9C  
Title : Crystal structure of YDCK from Salmonella cholerae. NESG target SCR6  
Authors : Benach, J.; Chen, Y.; Vorobiev, S.M.; Seetharaman, J.; Janjua, H.; Cooper, B.; Acton, X.T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2005-12-05  
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

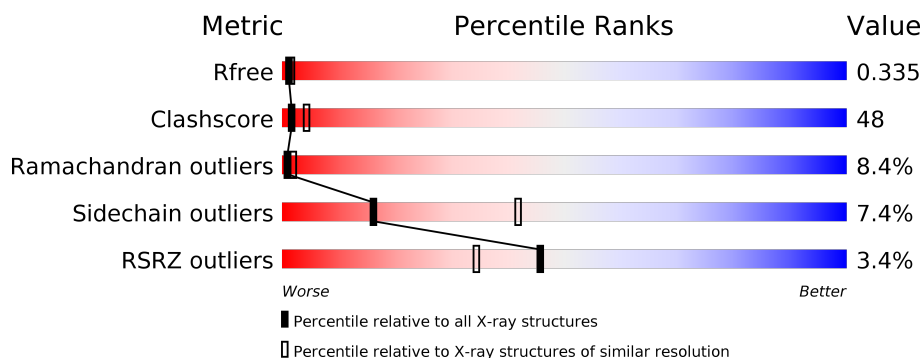
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	334	<div> <div>4%</div> <div> <div></div> <div>34%</div> <div>49%</div> <div>13%</div> <div>•</div> </div> </div>
1	B	334	<div> <div>2%</div> <div> <div></div> <div>31%</div> <div>52%</div> <div>11%</div> <div>• •</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5021 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 Hypothetical protein YDCK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2487	1546	452	481	8			
1	B	319	Total	C	N	O	S	0	0	0
			2478	1543	449	478	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	327	LEU	-	EXPRESSION TAG	GB 56413447
A	328	GLU	-	EXPRESSION TAG	GB 56413447
A	329	HIS	-	EXPRESSION TAG	GB 56413447
A	330	HIS	-	EXPRESSION TAG	GB 56413447
A	331	HIS	-	EXPRESSION TAG	GB 56413447
A	332	HIS	-	EXPRESSION TAG	GB 56413447
A	333	HIS	-	EXPRESSION TAG	GB 56413447
A	334	HIS	-	EXPRESSION TAG	GB 56413447
B	327	LEU	-	EXPRESSION TAG	GB 56413447
B	328	GLU	-	EXPRESSION TAG	GB 56413447
B	329	HIS	-	EXPRESSION TAG	GB 56413447
B	330	HIS	-	EXPRESSION TAG	GB 56413447
B	331	HIS	-	EXPRESSION TAG	GB 56413447
B	332	HIS	-	EXPRESSION TAG	GB 56413447
B	333	HIS	-	EXPRESSION TAG	GB 56413447
B	334	HIS	-	EXPRESSION TAG	GB 56413447

- Molecule 2 is CESIUM ION (three-letter code: CS) (formula: Cs).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Cs	0	0
			3	3		
2	A	3	Total	Cs	0	0
			3	3		

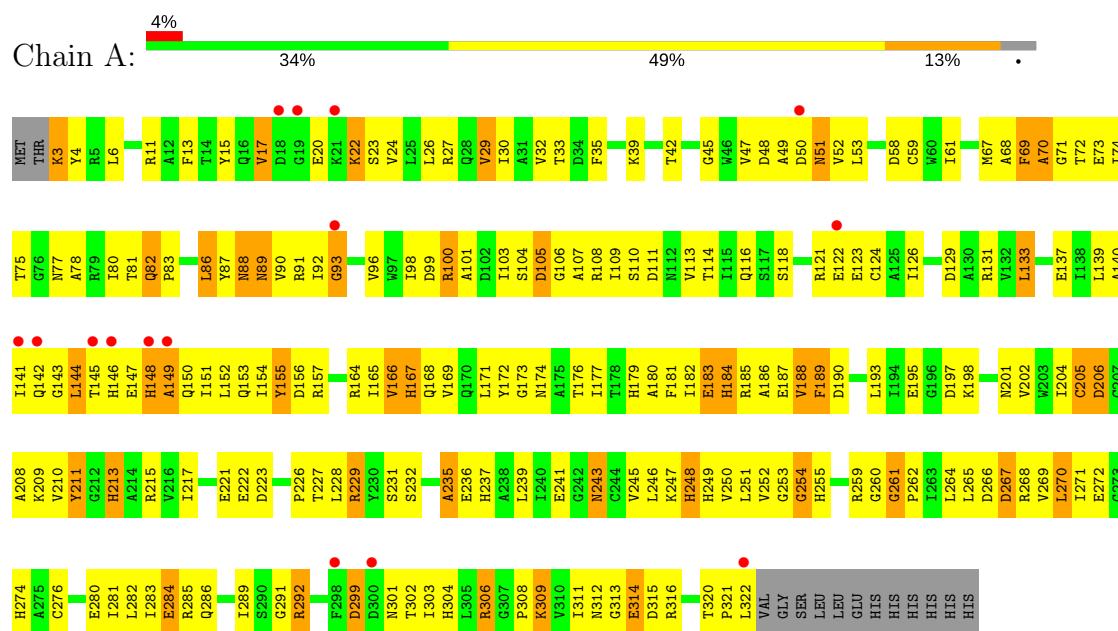
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total 25	O 25	0	0
3	B	25	Total 25	O 25	0	0

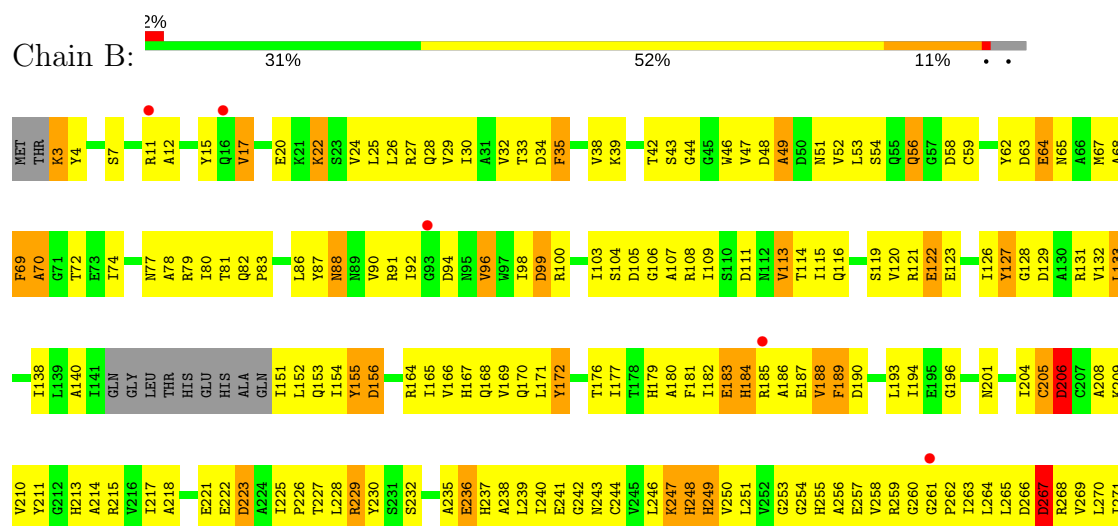
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Hypothetical protein YDCK



#### • Molecule 1: Hypothetical protein YDCK





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.60Å 104.60Å 162.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.99 – 2.70	Depositor EDS
% Data completeness (in resolution range)	78.1 (20.00-2.80) 89.4 (19.99-2.70)	Depositor EDS
$R_{merge}$	0.13	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.87 (at 2.71Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.253 , 0.323 0.276 , 0.335	Depositor DCC
$R_{free}$ test set	1966 reflections (9.81%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.8	Xtriage
Anisotropy	0.214	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 40.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	5021	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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

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	0/2529	0.88	10/3435 (0.3%)
1	B	0.38	0/2519	0.83	12/3420 (0.4%)
All	All	0.38	0/5048	0.86	22/6855 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	ALA	N-CA-C	6.59	128.78	111.00
1	B	69	PHE	N-CA-C	6.38	128.24	111.00
1	A	155	TYR	N-CA-C	6.31	128.05	111.00
1	A	189	PHE	N-CA-C	6.12	127.51	111.00
1	B	229	ARG	N-CA-C	6.05	127.35	111.00
1	A	213	HIS	N-CA-C	-5.87	95.14	111.00
1	B	172	TYR	N-CA-C	5.81	126.68	111.00
1	A	284	GLU	N-CA-C	5.72	126.44	111.00
1	A	229	ARG	N-CA-C	5.66	126.28	111.00
1	A	183	GLU	N-CA-C	5.56	126.01	111.00
1	A	254	GLY	N-CA-C	-5.53	99.29	113.10
1	B	284	GLU	N-CA-C	5.51	125.88	111.00
1	A	69	PHE	N-CA-C	5.50	125.86	111.00
1	B	127	TYR	N-CA-C	5.46	125.75	111.00
1	B	272	GLU	N-CA-C	5.43	125.67	111.00
1	A	205	CYS	N-CA-C	5.38	125.54	111.00
1	B	189	PHE	N-CA-C	5.34	125.41	111.00
1	B	155	TYR	N-CA-C	5.30	125.31	111.00
1	B	81	THR	N-CA-C	5.24	125.14	111.00
1	B	154	ILE	N-CA-C	-5.22	96.91	111.00
1	B	211	TYR	N-CA-C	5.19	125.02	111.00
1	B	183	GLU	N-CA-C	5.05	124.65	111.00



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	2487	0	2425	234	0
1	B	2478	0	2422	255	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	25	0	0	4	0
3	B	25	0	0	7	0
All	All	5021	0	4847	470	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 48.

All (470) 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:105:ASP:HB2	1:A:122:GLU:HA	1.40	1.03
1:A:96:VAL:HG23	1:A:110:SER:O	1.60	1.02
1:A:70:ALA:HB3	1:A:88:ASN:N	1.81	0.96
1:A:106:GLY:O	1:A:123:GLU:HA	1.68	0.93
1:B:268:ARG:HE	1:B:286:GLN:HE21	1.06	0.92
1:B:228:LEU:HD21	1:B:246:LEU:HD12	1.55	0.88
1:A:156:ASP:HB2	1:A:172:TYR:O	1.74	0.88
1:B:286:GLN:O	1:B:308:PRO:HA	1.75	0.86
1:A:144:LEU:H	1:A:144:LEU:HD22	1.42	0.85
1:A:114:THR:HG21	1:A:131:ARG:HH21	1.45	0.82
1:A:103:ILE:HG12	1:A:109:ILE:HD11	1.60	0.82
1:B:114:THR:HG21	1:B:131:ARG:HH21	1.45	0.81
1:B:121:ARG:HG3	1:B:122:GLU:H	1.45	0.80
1:B:87:TYR:O	1:B:88:ASN:HB2	1.81	0.80
1:A:291:GLY:HA3	1:A:312:ASN:O	1.81	0.80
1:A:70:ALA:HB3	1:A:88:ASN:H	1.43	0.80
1:A:106:GLY:HA3	1:A:123:GLU:HG2	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:ASP:H	1:B:77:ASN:ND2	1.79	0.80
1:B:121:ARG:HG3	1:B:122:GLU:N	1.98	0.79
1:A:39:LYS:O	1:A:42:THR:HG22	1.81	0.79
1:B:52:VAL:HG13	1:B:72:THR:O	1.82	0.79
1:B:169:VAL:HG23	1:B:183:GLU:O	1.82	0.79
1:A:147:GLU:HB2	1:A:185:ARG:NH2	1.98	0.78
1:B:58:ASP:H	1:B:77:ASN:HD22	1.28	0.78
1:B:28:GLN:HE22	1:B:64:GLU:HG2	1.47	0.77
1:B:268:ARG:HE	1:B:286:GLN:NE2	1.81	0.76
1:A:91:ARG:HB2	1:A:108:ARG:HG2	1.66	0.76
1:A:193:LEU:HD23	1:A:215:ARG:HG2	1.67	0.76
1:A:48:ASP:HB2	1:A:69:PHE:HA	1.67	0.75
1:B:106:GLY:HA3	1:B:123:GLU:HG2	1.66	0.75
1:A:313:GLY:O	1:A:315:ASP:N	2.17	0.75
1:A:292:ARG:CB	1:A:292:ARG:HH11	2.01	0.74
1:B:122:GLU:HB2	1:B:140:ALA:O	1.88	0.74
1:A:99:ASP:O	1:A:100:ARG:HB2	1.87	0.74
1:B:237:HIS:HB2	1:B:254:GLY:O	1.87	0.73
1:A:274:HIS:HB3	1:B:131:ARG:HD2	1.71	0.73
1:B:166:VAL:HG12	1:B:167:HIS:H	1.53	0.72
1:B:169:VAL:HG22	1:B:186:ALA:HB3	1.72	0.72
1:A:152:LEU:HD12	1:A:166:VAL:O	1.90	0.72
1:B:59:CYS:HA	1:B:78:ALA:O	1.90	0.72
1:A:177:ILE:HD11	1:A:188:VAL:HG21	1.73	0.71
1:B:4:TYR:HA	1:B:32:VAL:HG23	1.71	0.71
1:A:75:THR:O	1:A:93:GLY:HA2	1.91	0.71
1:B:171:LEU:HD23	1:B:188:VAL:HG13	1.72	0.71
1:B:72:THR:HG23	1:B:90:VAL:HB	1.71	0.71
1:A:182:ILE:HG12	1:A:204:ILE:HD12	1.72	0.71
1:A:215:ARG:HD2	1:B:193:LEU:HD13	1.73	0.70
1:A:285:ARG:CZ	1:A:308:PRO:HD3	2.21	0.70
1:B:96:VAL:HG21	1:B:109:ILE:HG22	1.72	0.70
1:A:58:ASP:H	1:A:77:ASN:ND2	1.90	0.69
1:A:3:LYS:HB3	1:A:33:THR:CG2	2.23	0.69
1:B:170:GLN:HB2	1:B:187:GLU:HG2	1.74	0.69
1:A:272:GLU:HG2	1:A:272:GLU:O	1.93	0.69
1:B:236:GLU:HB2	1:B:254:GLY:N	2.08	0.68
1:B:121:ARG:CG	1:B:122:GLU:H	2.00	0.68
1:A:262:PRO:HD2	1:A:280:GLU:HB2	1.74	0.68
1:B:236:GLU:HB2	1:B:254:GLY:H	1.59	0.68
1:B:82:GLN:HB3	1:B:83:PRO:CD	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:228:LEU:HD23	1:B:246:LEU:HB2	1.75	0.67
1:A:251:LEU:C	1:A:251:LEU:HD23	2.14	0.67
1:A:3:LYS:HB3	1:A:33:THR:HG22	1.77	0.67
1:A:82:GLN:HB3	1:A:83:PRO:CD	2.25	0.67
1:A:267:ASP:HB2	1:A:284:GLU:O	1.95	0.67
1:A:292:ARG:HB2	1:A:292:ARG:HH11	1.59	0.67
1:A:241:GLU:OE2	1:A:259:ARG:HB2	1.94	0.67
1:A:301:ASN:HA	3:A:421:HOH:O	1.95	0.67
1:A:147:GLU:HB2	1:A:185:ARG:HH21	1.59	0.66
1:A:166:VAL:HG23	1:A:183:GLU:HG2	1.77	0.66
1:A:48:ASP:HB2	1:A:69:PHE:CA	2.23	0.66
1:B:53:LEU:HD23	1:B:54:SER:N	2.11	0.66
1:A:271:ILE:HD12	1:A:271:ILE:N	2.11	0.65
1:A:22:LYS:HE2	1:A:23:SER:H	1.60	0.65
1:A:169:VAL:HG21	1:A:182:ILE:HG22	1.78	0.65
1:B:114:THR:HG21	1:B:131:ARG:NH2	2.11	0.65
1:A:59:CYS:HA	1:A:78:ALA:O	1.95	0.65
1:B:115:ILE:HD13	1:B:132:VAL:HB	1.79	0.65
1:A:116:GLN:HE22	1:B:296:ILE:HD11	1.62	0.65
1:A:316:ARG:HD2	1:B:99:ASP:OD1	1.96	0.64
1:B:90:VAL:O	1:B:91:ARG:HD2	1.97	0.64
1:B:17:VAL:CG2	1:B:20:GLU:HB2	2.28	0.64
1:B:155:TYR:CD1	1:B:155:TYR:O	2.51	0.64
1:B:166:VAL:HG12	1:B:167:HIS:N	2.12	0.64
1:A:285:ARG:NH2	1:A:308:PRO:HD3	2.13	0.64
1:B:187:GLU:O	1:B:209:LYS:HA	1.97	0.64
1:B:56:GLN:N	1:B:56:GLN:NE2	2.46	0.64
1:A:169:VAL:HG23	1:A:183:GLU:O	1.97	0.64
1:A:171:LEU:HD23	1:A:188:VAL:HG13	1.80	0.64
1:B:111:ASP:HB2	1:B:127:TYR:O	1.98	0.64
1:A:11:ARG:HD2	1:A:13:PHE:CZ	2.33	0.64
1:A:100:ARG:O	1:A:100:ARG:HD2	1.97	0.63
1:A:260:GLY:O	1:A:261:GLY:O	2.16	0.63
1:A:148:HIS:O	1:A:150:GLN:N	2.30	0.63
1:A:131:ARG:HH11	1:B:274:HIS:HB3	1.64	0.62
1:A:166:VAL:CG2	1:A:183:GLU:HG2	2.28	0.62
1:B:281:ILE:N	1:B:281:ILE:HD12	2.14	0.62
1:B:260:GLY:HA3	1:B:279:GLY:H	1.64	0.62
1:B:270:LEU:CD2	1:B:272:GLU:HB2	2.28	0.62
1:A:155:TYR:O	1:A:155:TYR:HD1	1.83	0.62
1:A:52:VAL:HG22	1:A:72:THR:O	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:313:GLY:C	1:B:315:ASP:H	2.02	0.62
1:B:312:ASN:O	1:B:313:GLY:O	2.17	0.61
1:A:101:ALA:HB1	1:A:118:SER:O	2.00	0.61
1:B:306:ARG:H	1:B:326:LEU:CD2	2.14	0.61
1:B:189:PHE:O	1:B:189:PHE:HD1	1.84	0.61
1:B:103:ILE:N	1:B:103:ILE:HD12	2.14	0.61
1:B:70:ALA:HB3	1:B:88:ASN:H	1.65	0.61
1:A:86:LEU:N	1:A:86:LEU:HD23	2.15	0.61
1:A:205:CYS:HB3	3:A:417:HOH:O	2.01	0.60
1:B:176:THR:HB	1:B:193:LEU:HD12	1.82	0.60
1:B:306:ARG:H	1:B:326:LEU:HD21	1.66	0.60
1:A:29:VAL:HB	1:A:45:GLY:O	2.01	0.60
1:A:155:TYR:O	1:A:155:TYR:CD1	2.55	0.60
1:B:103:ILE:CG2	1:B:107:ALA:HB3	2.32	0.60
1:A:99:ASP:OD1	1:B:316:ARG:HD2	2.02	0.60
1:A:86:LEU:HA	1:A:103:ILE:O	2.02	0.60
1:A:122:GLU:HG3	1:A:141:ILE:HD13	1.83	0.60
1:A:87:TYR:O	1:A:88:ASN:HB2	2.02	0.60
1:A:270:LEU:HD21	1:A:272:GLU:HB2	1.84	0.60
1:B:171:LEU:CD2	1:B:188:VAL:HG13	2.32	0.60
1:B:313:GLY:O	1:B:315:ASP:N	2.30	0.60
1:A:58:ASP:H	1:A:77:ASN:HD22	1.50	0.59
1:A:50:ASP:O	1:A:51:ASN:HB2	2.00	0.59
1:A:73:GLU:HG3	1:A:91:ARG:NH1	2.18	0.59
1:B:306:ARG:HB2	1:B:326:LEU:HD22	1.85	0.59
1:A:285:ARG:HG3	1:A:306:ARG:CZ	2.33	0.59
1:B:177:ILE:N	1:B:177:ILE:HD12	2.18	0.59
1:A:152:LEU:C	1:A:152:LEU:HD23	2.23	0.59
1:A:70:ALA:HB3	1:A:88:ASN:CA	2.33	0.59
1:B:227:THR:HG22	1:B:229:ARG:HG3	1.85	0.59
1:A:206:ASP:OD2	1:A:229:ARG:O	2.20	0.58
1:A:313:GLY:C	1:A:315:ASP:H	2.04	0.58
1:B:56:GLN:NE2	1:B:56:GLN:H	2.02	0.58
1:B:167:HIS:HB3	1:B:168:GLN:OE1	2.03	0.58
1:B:44:GLY:O	1:B:64:GLU:HA	2.03	0.58
1:A:157:ARG:HH11	1:A:157:ARG:HG3	1.67	0.58
1:B:261:GLY:O	1:B:263:ILE:HG13	2.03	0.58
1:B:7:SER:HB2	1:B:28:GLN:HB3	1.85	0.58
1:B:109:ILE:HD13	1:B:115:ILE:HG13	1.85	0.57
1:B:152:LEU:HD23	1:B:153:GLN:N	2.19	0.57
1:A:253:GLY:O	1:A:272:GLU:HG3	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:ALA:HB3	1:B:88:ASN:N	2.19	0.57
1:B:4:TYR:CD1	1:B:29:VAL:HG13	2.39	0.57
1:B:304:HIS:HD2	1:B:305:LEU:N	2.03	0.57
1:B:103:ILE:HG23	1:B:107:ALA:HB3	1.86	0.57
1:B:86:LEU:CD2	1:B:103:ILE:HD13	2.35	0.56
1:A:153:GLN:HB3	1:A:155:TYR:HD2	1.69	0.56
1:A:264:LEU:HB3	1:A:282:LEU:HD12	1.88	0.56
1:B:47:VAL:HG12	1:B:49:ALA:H	1.70	0.56
1:A:217:ILE:O	1:A:226:PRO:HG2	2.06	0.56
1:A:187:GLU:O	1:A:209:LYS:HA	2.06	0.56
1:B:217:ILE:O	1:B:226:PRO:HG2	2.05	0.56
1:B:68:ALA:HA	1:B:86:LEU:O	2.06	0.56
1:A:90:VAL:O	1:A:91:ARG:HD2	2.04	0.56
1:B:206:ASP:CG	1:B:229:ARG:O	2.44	0.56
1:B:63:ASP:OD1	1:B:65:ASN:HB2	2.06	0.56
1:B:82:GLN:HB3	1:B:83:PRO:HD2	1.87	0.55
1:B:187:GLU:HB2	1:B:209:LYS:HG2	1.88	0.55
1:A:285:ARG:HH21	1:A:286:GLN:NE2	2.04	0.55
1:A:70:ALA:HB3	1:A:88:ASN:HA	1.88	0.55
1:B:86:LEU:HD22	1:B:103:ILE:HB	1.88	0.55
1:A:247:LYS:HB2	1:A:266:ASP:HA	1.88	0.55
1:B:47:VAL:CG1	1:B:48:ASP:N	2.69	0.55
1:A:103:ILE:N	1:A:103:ILE:HD12	2.22	0.55
1:B:129:ASP:N	1:B:156:ASP:O	2.35	0.55
1:A:221:GLU:H	1:A:221:GLU:CD	2.10	0.55
1:A:281:ILE:HG22	1:A:283:ILE:HG13	1.86	0.55
1:B:166:VAL:O	1:B:183:GLU:O	2.25	0.55
1:B:256:ALA:CB	1:B:271:ILE:HG22	2.37	0.55
1:A:13:PHE:O	1:A:23:SER:HA	2.07	0.55
1:B:168:GLN:HE21	1:B:185:ARG:NH1	2.05	0.55
1:B:222:GLU:HG2	1:B:223:ASP:OD1	2.07	0.55
1:B:268:ARG:NE	1:B:286:GLN:HE21	1.90	0.54
1:A:169:VAL:CG2	1:A:182:ILE:HG22	2.38	0.54
1:A:270:LEU:CD2	1:A:272:GLU:HB2	2.38	0.54
1:B:96:VAL:CG2	1:B:109:ILE:HG22	2.35	0.54
1:A:70:ALA:CB	1:A:88:ASN:HA	2.37	0.54
1:B:214:ALA:HB2	1:B:235:ALA:O	2.08	0.54
1:B:268:ARG:HH11	1:B:268:ARG:HG3	1.71	0.54
1:B:168:GLN:HE21	1:B:185:ARG:HH12	1.56	0.54
1:A:68:ALA:HA	1:A:86:LEU:O	2.08	0.54
1:B:189:PHE:CD1	1:B:189:PHE:O	2.60	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:THR:HB	1:A:245:VAL:HG22	1.88	0.54
1:A:116:GLN:NE2	1:B:296:ILE:HD11	2.23	0.54
1:B:86:LEU:HD21	1:B:103:ILE:HD13	1.89	0.54
1:A:91:ARG:CB	1:A:108:ARG:HG2	2.37	0.54
1:B:239:LEU:HB3	1:B:257:GLU:HG3	1.90	0.54
1:B:290:SER:OG	1:B:290:SER:O	2.22	0.54
1:A:177:ILE:HD11	1:A:188:VAL:CG2	2.38	0.53
1:A:272:GLU:CG	1:A:272:GLU:O	2.56	0.53
1:B:228:LEU:CD2	1:B:246:LEU:HD12	2.33	0.53
1:A:157:ARG:NH1	1:A:157:ARG:HG3	2.23	0.53
1:A:193:LEU:CD2	1:A:215:ARG:HG2	2.37	0.53
1:A:285:ARG:HB2	1:A:306:ARG:O	2.09	0.53
1:A:281:ILE:N	1:A:281:ILE:HD12	2.23	0.53
1:B:152:LEU:HD23	1:B:152:LEU:C	2.29	0.53
1:A:144:LEU:H	1:A:144:LEU:CD2	2.17	0.53
1:A:3:LYS:HB2	1:A:3:LYS:NZ	2.23	0.53
1:A:148:HIS:O	1:A:149:ALA:C	2.47	0.53
1:A:105:ASP:N	1:A:121:ARG:O	2.41	0.53
1:A:172:TYR:CE1	1:A:189:PHE:HB2	2.43	0.53
1:A:87:TYR:O	1:A:88:ASN:CB	2.56	0.53
1:B:87:TYR:O	1:B:104:SER:O	2.26	0.53
1:B:11:ARG:HG2	1:B:12:ALA:H	1.73	0.52
1:B:229:ARG:NH1	1:B:247:LYS:HE2	2.24	0.52
1:B:268:ARG:NH1	1:B:268:ARG:HG3	2.24	0.52
1:B:32:VAL:HG12	1:B:32:VAL:O	2.09	0.52
1:B:91:ARG:HB2	1:B:108:ARG:HG2	1.90	0.52
1:B:108:ARG:HG3	1:B:108:ARG:HH11	1.74	0.52
1:B:230:TYR:C	1:B:232:SER:H	2.12	0.52
1:B:267:ASP:HB2	1:B:284:GLU:O	2.09	0.52
1:A:151:ILE:O	1:A:153:GLN:HG3	2.08	0.52
1:A:88:ASN:HB2	1:A:104:SER:O	2.09	0.52
1:A:270:LEU:C	1:A:271:ILE:HD12	2.30	0.52
1:B:306:ARG:CB	1:B:326:LEU:HD22	2.39	0.52
1:A:241:GLU:O	1:A:259:ARG:HA	2.09	0.52
1:B:120:VAL:HA	1:B:138:ILE:HB	1.92	0.52
1:A:129:ASP:HB3	1:B:274:HIS:HE1	1.75	0.52
1:A:82:GLN:CB	1:A:83:PRO:CD	2.88	0.52
1:B:269:VAL:HG12	1:B:271:ILE:HD12	1.92	0.52
1:B:27:ARG:NE	3:B:408:HOH:O	2.42	0.52
1:A:243:ASN:HD22	1:A:243:ASN:C	2.14	0.52
1:B:312:ASN:ND2	1:B:313:GLY:H	2.08	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:GLY:HA3	3:B:425:HOH:O	2.09	0.51
1:B:299:ASP:C	1:B:301:ASN:N	2.63	0.51
1:A:47:VAL:CG1	1:A:52:VAL:HB	2.39	0.51
1:B:164:ARG:C	1:B:165:ILE:HD13	2.30	0.51
1:B:215:ARG:HB2	1:B:239:LEU:HD23	1.92	0.51
1:B:256:ALA:HB1	1:B:271:ILE:HG22	1.91	0.51
1:B:291:GLY:HA3	1:B:312:ASN:O	2.10	0.51
1:B:70:ALA:HB3	1:B:88:ASN:CA	2.41	0.51
1:A:248:HIS:HB3	1:A:249:HIS:HD2	1.75	0.51
1:B:35:PHE:CD1	1:B:35:PHE:N	2.79	0.51
1:B:28:GLN:NE2	1:B:43:SER:HB2	2.24	0.51
1:A:98:ILE:HD13	1:A:103:ILE:HD13	1.93	0.51
1:A:89:ASN:O	1:A:91:ARG:HD3	2.11	0.50
1:A:215:ARG:HD2	1:B:193:LEU:CD1	2.41	0.50
1:B:88:ASN:ND2	1:B:105:ASP:OD1	2.44	0.50
1:A:195:GLU:CD	1:B:259:ARG:HH22	2.14	0.50
1:B:303:ILE:HD11	1:B:318:THR:HA	1.93	0.50
1:A:303:ILE:HA	1:A:320:THR:OG1	2.12	0.50
1:B:156:ASP:OD2	1:B:172:TYR:O	2.29	0.50
1:B:272:GLU:HB3	1:B:290:SER:HB2	1.93	0.50
1:A:284:GLU:OE1	1:A:306:ARG:HD2	2.11	0.50
1:A:140:ALA:HB2	1:A:152:LEU:HD13	1.94	0.50
1:A:167:HIS:HB2	1:A:184:HIS:HA	1.94	0.50
1:A:48:ASP:CB	1:A:69:PHE:HA	2.39	0.50
1:A:264:LEU:O	1:A:265:LEU:HD23	2.11	0.50
1:B:133:LEU:HD23	1:B:133:LEU:N	2.26	0.50
1:A:131:ARG:HD2	1:B:274:HIS:HB3	1.93	0.50
1:A:113:VAL:CG1	1:A:114:THR:N	2.75	0.50
1:A:74:ILE:HD13	1:A:80:ILE:CD1	2.42	0.50
1:B:22:LYS:HE2	1:B:22:LYS:HA	1.93	0.50
1:B:305:LEU:HD13	1:B:311:ILE:CD1	2.42	0.50
1:A:270:LEU:HD23	1:A:271:ILE:N	2.27	0.50
1:A:82:GLN:HB3	1:A:83:PRO:HD2	1.94	0.49
1:B:100:ARG:O	1:B:100:ARG:HG2	2.12	0.49
1:A:237:HIS:HB2	1:A:254:GLY:O	2.12	0.49
1:A:49:ALA:O	1:A:50:ASP:C	2.50	0.49
1:B:184:HIS:CD2	1:B:184:HIS:N	2.80	0.49
1:A:167:HIS:HB3	1:A:168:GLN:OE1	2.12	0.49
1:A:271:ILE:HA	1:A:289:ILE:O	2.12	0.49
1:A:4:TYR:HB2	1:A:30:ILE:O	2.13	0.49
1:B:103:ILE:N	1:B:103:ILE:CD1	2.74	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:ILE:HG22	1:B:283:ILE:HG13	1.94	0.49
1:B:267:ASP:HB2	1:B:285:ARG:HA	1.95	0.49
1:A:145:THR:O	1:A:145:THR:HG22	2.13	0.49
1:A:280:GLU:C	1:A:281:ILE:HD12	2.33	0.49
1:B:123:GLU:C	1:B:151:ILE:HD11	2.33	0.49
1:B:285:ARG:NH1	3:B:419:HOH:O	2.45	0.49
1:B:86:LEU:HD23	1:B:103:ILE:H	1.78	0.49
1:A:103:ILE:CG2	1:A:107:ALA:HB3	2.42	0.48
1:A:156:ASP:CB	1:A:172:TYR:O	2.56	0.48
1:A:285:ARG:HG3	1:A:306:ARG:NH2	2.28	0.48
1:B:90:VAL:HG13	1:B:103:ILE:HG21	1.94	0.48
1:B:53:LEU:HD23	1:B:54:SER:H	1.78	0.48
1:B:28:GLN:CD	1:B:43:SER:HB2	2.32	0.48
1:A:180:ALA:C	1:A:181:PHE:HD2	2.16	0.48
1:B:194:ILE:CD1	1:B:204:ILE:HD11	2.43	0.48
1:A:124:CYS:HA	1:A:151:ILE:CG1	2.44	0.48
1:A:137:GLU:CD	1:A:164:ARG:HH21	2.17	0.48
1:A:144:LEU:N	1:A:144:LEU:HD22	2.21	0.48
1:A:92:ILE:HG23	1:A:96:VAL:HG11	1.95	0.48
1:A:172:TYR:CD1	1:A:189:PHE:HB2	2.49	0.48
1:A:282:LEU:HD23	1:A:304:HIS:CD2	2.48	0.48
1:B:188:VAL:HA	1:B:210:VAL:O	2.14	0.48
1:A:285:ARG:NH2	1:A:286:GLN:NE2	2.62	0.48
1:B:4:TYR:HB2	1:B:30:ILE:O	2.14	0.48
1:B:87:TYR:O	1:B:88:ASN:CB	2.59	0.48
1:B:168:GLN:NE2	1:B:185:ARG:NH1	2.62	0.48
1:B:3:LYS:NZ	1:B:3:LYS:HB2	2.29	0.48
1:A:264:LEU:C	1:A:265:LEU:HD23	2.34	0.48
1:B:182:ILE:HA	1:B:204:ILE:O	2.14	0.47
1:A:122:GLU:HB2	1:A:140:ALA:O	2.14	0.47
1:A:292:ARG:HB2	1:A:292:ARG:NH1	2.28	0.47
1:B:91:ARG:CB	1:B:108:ARG:HG2	2.44	0.47
1:A:227:THR:HG22	1:A:229:ARG:HG3	1.96	0.47
1:A:92:ILE:HG12	1:A:109:ILE:HD12	1.97	0.47
1:A:235:ALA:O	1:A:236:GLU:HG3	2.14	0.47
1:A:47:VAL:HG12	1:A:48:ASP:N	2.28	0.47
1:A:81:THR:HB	1:A:82:GLN:HE21	1.78	0.47
1:B:182:ILE:HG22	1:B:186:ALA:HB3	1.96	0.47
1:B:184:HIS:HB2	1:B:206:ASP:O	2.14	0.47
1:B:249:HIS:CD2	3:B:409:HOH:O	2.67	0.47
1:B:237:HIS:CB	1:B:254:GLY:O	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:GLU:CD	1:B:314:GLU:H	2.18	0.47
1:B:92:ILE:HD13	1:B:98:ILE:HD12	1.95	0.47
1:A:153:GLN:HB3	1:A:155:TYR:CD2	2.48	0.47
1:A:186:ALA:HB1	1:A:204:ILE:HG22	1.96	0.47
1:A:321:PRO:O	1:A:322:LEU:HG	2.15	0.47
1:B:38:VAL:HG13	1:B:42:THR:HG21	1.97	0.47
1:A:267:ASP:HB3	1:A:268:ARG:H	1.45	0.47
1:B:240:ILE:HD13	1:B:258:VAL:HB	1.97	0.47
1:B:266:ASP:O	1:B:269:VAL:HG23	2.14	0.47
1:B:47:VAL:HG12	1:B:49:ALA:N	2.29	0.47
1:A:104:SER:HB2	1:A:121:ARG:HA	1.96	0.47
1:A:70:ALA:CB	1:A:88:ASN:N	2.68	0.47
1:B:246:LEU:HD22	1:B:265:LEU:HD12	1.96	0.47
1:B:90:VAL:HG22	1:B:103:ILE:HG22	1.95	0.47
1:B:217:ILE:HG22	1:B:218:ALA:N	2.30	0.46
1:B:25:LEU:O	1:B:26:LEU:HD23	2.15	0.46
1:A:6:LEU:HD22	1:A:27:ARG:HB2	1.97	0.46
1:B:266:ASP:O	1:B:267:ASP:C	2.53	0.46
1:B:74:ILE:HD13	1:B:80:ILE:CD1	2.46	0.46
1:A:171:LEU:CD2	1:A:188:VAL:HG13	2.45	0.46
1:B:312:ASN:ND2	1:B:313:GLY:N	2.64	0.46
1:B:69:PHE:N	1:B:69:PHE:CD2	2.84	0.46
1:B:69:PHE:CZ	1:B:87:TYR:HB3	2.50	0.46
1:A:195:GLU:OE1	1:B:259:ARG:NH2	2.46	0.46
1:B:63:ASP:OD1	1:B:65:ASN:N	2.49	0.46
1:A:198:LYS:NZ	1:B:221:GLU:HG3	2.31	0.46
1:A:3:LYS:HZ3	1:A:3:LYS:HB2	1.81	0.46
1:B:113:VAL:CG1	1:B:114:THR:N	2.78	0.46
1:B:169:VAL:HG21	1:B:182:ILE:HG22	1.98	0.46
1:A:131:ARG:HB3	1:A:133:LEU:HD21	1.97	0.45
1:A:281:ILE:CG2	1:A:283:ILE:HG13	2.45	0.45
1:A:289:ILE:HA	1:A:311:ILE:O	2.15	0.45
1:A:316:ARG:HH11	1:A:316:ARG:HG2	1.81	0.45
1:B:289:ILE:HA	1:B:311:ILE:O	2.15	0.45
1:A:213:HIS:HE1	1:B:213:HIS:CE1	2.34	0.45
1:B:299:ASP:C	1:B:301:ASN:H	2.19	0.45
1:B:58:ASP:N	1:B:77:ASN:ND2	2.57	0.45
1:A:179:HIS:O	1:A:202:VAL:HG23	2.16	0.45
1:A:98:ILE:HD13	1:A:103:ILE:CD1	2.47	0.45
1:A:314:GLU:HG3	1:B:79:ARG:HE	1.82	0.45
1:A:72:THR:HG23	1:A:90:VAL:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:ARG:HG2	1:B:181:PHE:CD2	2.51	0.45
1:B:15:TYR:HE2	1:B:24:VAL:HG13	1.81	0.45
1:A:157:ARG:NH1	3:A:420:HOH:O	2.47	0.45
1:A:246:LEU:HD23	1:A:246:LEU:N	2.31	0.45
1:A:313:GLY:C	1:A:315:ASP:N	2.65	0.45
1:A:232:SER:HA	1:A:250:VAL:O	2.17	0.45
1:A:87:TYR:O	1:A:104:SER:O	2.35	0.45
1:A:274:HIS:CB	1:B:131:ARG:HD2	2.44	0.45
1:B:281:ILE:N	1:B:281:ILE:CD1	2.80	0.45
1:A:15:TYR:CE2	1:A:22:LYS:HB3	2.52	0.45
1:B:313:GLY:C	1:B:315:ASP:N	2.70	0.45
1:B:304:HIS:CD2	1:B:305:LEU:N	2.84	0.45
1:A:174:ASN:OD1	1:B:255:HIS:CE1	2.70	0.44
1:B:28:GLN:NE2	1:B:64:GLU:HG2	2.25	0.44
1:B:292:ARG:HH12	1:B:314:GLU:HB3	1.82	0.44
1:B:38:VAL:HB	3:B:407:HOH:O	2.18	0.44
1:A:103:ILE:CG1	1:A:109:ILE:HD11	2.39	0.44
1:B:63:ASP:C	1:B:65:ASN:H	2.21	0.44
1:A:251:LEU:HD23	1:A:252:VAL:N	2.32	0.44
1:A:6:LEU:HD11	1:A:50:ASP:H	1.82	0.44
1:A:74:ILE:HD13	1:A:80:ILE:HD12	2.00	0.44
1:B:285:ARG:HG3	1:B:306:ARG:NH2	2.32	0.44
1:B:33:THR:HG23	1:B:34:ASP:N	2.33	0.44
1:B:151:ILE:N	3:B:410:HOH:O	2.50	0.44
1:B:67:MET:HB3	1:B:69:PHE:CE2	2.53	0.44
1:A:188:VAL:HA	1:A:210:VAL:O	2.17	0.44
1:B:113:VAL:HG11	1:B:126:ILE:CG2	2.48	0.44
1:B:312:ASN:CG	1:B:313:GLY:H	2.22	0.44
1:A:316:ARG:HG2	1:A:316:ARG:NH1	2.33	0.44
1:A:71:GLY:O	1:A:91:ARG:NH1	2.44	0.44
1:B:249:HIS:CD2	1:B:249:HIS:H	2.35	0.44
1:B:306:ARG:HH11	1:B:306:ARG:HG2	1.83	0.44
1:A:88:ASN:C	1:A:90:VAL:N	2.71	0.43
1:B:256:ALA:HB1	1:B:271:ILE:CG2	2.48	0.43
1:A:174:ASN:N	1:A:190:ASP:O	2.45	0.43
1:B:249:HIS:N	1:B:249:HIS:CD2	2.86	0.43
1:A:271:ILE:CD1	1:A:271:ILE:N	2.81	0.43
1:A:4:TYR:HA	1:A:32:VAL:HG23	2.00	0.43
1:B:165:ILE:HG22	1:B:169:VAL:HB	2.00	0.43
1:B:166:VAL:CG1	1:B:167:HIS:H	2.27	0.43
1:A:149:ALA:O	1:A:151:ILE:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:LYS:HE3	1:A:264:LEU:HD11	2.00	0.43
1:A:17:VAL:CG2	1:A:20:GLU:HB2	2.49	0.43
1:A:48:ASP:OD2	1:A:69:PHE:O	2.37	0.43
1:B:47:VAL:HG12	1:B:48:ASP:N	2.33	0.43
1:A:309:LYS:NZ	1:A:309:LYS:HB2	2.33	0.43
1:B:171:LEU:CD2	1:B:188:VAL:CG1	2.96	0.43
1:A:126:ILE:HG23	1:A:154:ILE:HB	2.00	0.43
1:B:238:ALA:HB2	1:B:253:GLY:HA2	2.01	0.43
1:B:74:ILE:HD13	1:B:80:ILE:HD12	2.00	0.43
1:B:240:ILE:HA	1:B:258:VAL:O	2.19	0.43
1:B:90:VAL:HG12	1:B:91:ARG:N	2.33	0.43
1:A:261:GLY:N	3:A:409:HOH:O	2.52	0.43
1:A:61:ILE:HA	1:A:80:ILE:HB	2.01	0.43
1:B:48:ASP:HB2	1:B:69:PHE:CA	2.49	0.43
1:A:190:ASP:OD2	1:A:211:TYR:O	2.37	0.42
1:B:206:ASP:HB2	1:B:230:TYR:HA	2.01	0.42
1:B:248:HIS:O	1:B:250:VAL:N	2.52	0.42
1:B:280:GLU:O	1:B:280:GLU:HG2	2.19	0.42
1:A:106:GLY:C	1:A:123:GLU:HA	2.38	0.42
1:B:155:TYR:O	1:B:156:ASP:CG	2.58	0.42
1:B:153:GLN:HB2	1:B:170:GLN:NE2	2.34	0.42
1:B:306:ARG:HH11	1:B:306:ARG:CG	2.32	0.42
1:A:228:LEU:HD23	1:A:228:LEU:N	2.34	0.42
1:B:113:VAL:HG12	1:B:114:THR:N	2.34	0.42
1:B:48:ASP:CB	1:B:69:PHE:O	2.68	0.42
1:A:103:ILE:HG23	1:A:107:ALA:HB3	2.01	0.42
1:A:121:ARG:HB3	1:A:139:LEU:HD12	2.02	0.42
1:A:140:ALA:HB2	1:A:152:LEU:CD1	2.50	0.42
1:B:179:HIS:O	1:B:196:GLY:HA3	2.20	0.42
1:B:321:PRO:O	1:B:326:LEU:HG	2.20	0.42
1:B:248:HIS:HB3	1:B:249:HIS:H	1.59	0.42
1:A:182:ILE:HA	1:A:204:ILE:O	2.20	0.42
1:A:215:ARG:CD	1:B:193:LEU:HD22	2.49	0.42
1:A:222:GLU:HG2	1:A:223:ASP:OD1	2.20	0.42
1:A:237:HIS:CD2	1:A:255:HIS:CE1	3.08	0.42
1:B:169:VAL:CG2	1:B:182:ILE:HG22	2.50	0.42
1:B:184:HIS:CD2	1:B:184:HIS:H	2.37	0.42
1:B:28:GLN:N	1:B:46:TRP:CE3	2.88	0.42
1:A:90:VAL:HG13	1:A:103:ILE:HG21	2.02	0.41
1:B:109:ILE:CD1	1:B:115:ILE:HG13	2.49	0.41
1:B:264:LEU:HD23	1:B:282:LEU:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:LEU:HD23	1:A:153:GLN:N	2.36	0.41
1:B:165:ILE:HD13	1:B:165:ILE:N	2.34	0.41
1:B:306:ARG:NE	3:B:420:HOH:O	2.52	0.41
1:A:165:ILE:N	1:A:165:ILE:HD13	2.35	0.41
1:A:124:CYS:HA	1:A:151:ILE:HG12	2.02	0.41
1:A:285:ARG:HH21	1:A:286:GLN:CD	2.24	0.41
1:B:108:ARG:NH1	1:B:108:ARG:HG3	2.36	0.41
1:B:285:ARG:NH2	1:B:286:GLN:OE1	2.53	0.41
1:B:299:ASP:HB3	1:B:300:ASP:H	1.66	0.41
1:B:39:LYS:HA	1:B:39:LYS:HD2	1.89	0.41
1:A:88:ASN:O	1:A:90:VAL:N	2.54	0.41
1:B:116:GLN:O	1:B:133:LEU:HA	2.19	0.41
1:B:166:VAL:HB	1:B:167:HIS:ND1	2.36	0.41
1:A:266:ASP:O	1:A:269:VAL:HG23	2.19	0.41
1:B:194:ILE:HD11	1:B:204:ILE:HD11	2.02	0.41
1:B:226:PRO:HA	1:B:244:CYS:O	2.21	0.41
1:B:104:SER:HB2	1:B:121:ARG:HA	2.01	0.41
1:A:215:ARG:CD	1:B:193:LEU:HD13	2.46	0.41
1:A:180:ALA:CB	1:A:202:VAL:HB	2.51	0.41
1:A:284:GLU:OE1	1:A:304:HIS:NE2	2.54	0.41
1:B:62:TYR:CE1	1:B:79:ARG:HG2	2.55	0.41
1:B:86:LEU:HD13	1:B:92:ILE:CD1	2.51	0.41
1:A:122:GLU:CG	1:A:141:ILE:HD13	2.51	0.41
1:A:190:ASP:HB2	1:A:211:TYR:O	2.21	0.41
1:B:267:ASP:HB3	1:B:268:ARG:H	1.51	0.41
1:B:4:TYR:OH	1:B:53:LEU:O	2.39	0.41
1:B:201:ASN:O	1:B:225:ILE:HG23	2.21	0.41
1:B:304:HIS:C	1:B:304:HIS:CD2	2.94	0.41
1:A:215:ARG:HB2	1:A:239:LEU:HD22	2.04	0.40
1:A:48:ASP:HB2	1:A:69:PHE:CB	2.51	0.40
1:A:26:LEU:CD1	1:A:67:MET:HG2	2.50	0.40
1:B:241:GLU:CD	1:B:242:GLY:N	2.75	0.40
1:A:179:HIS:ND1	1:A:197:ASP:HA	2.37	0.40
1:B:205:CYS:HB2	1:B:206:ASP:H	1.63	0.40
1:B:208:ALA:HA	1:B:232:SER:O	2.21	0.40
1:B:269:VAL:HG12	1:B:270:LEU:N	2.35	0.40
1:B:267:ASP:CB	1:B:285:ARG:HA	2.52	0.40
1:B:48:ASP:O	1:B:49:ALA:HB2	2.22	0.40
1:A:168:GLN:HE21	1:A:185:ARG:NH1	2.18	0.40
1:A:201:ASN:HB3	1:A:202:VAL:H	1.77	0.40
1:A:47:VAL:HG11	1:A:52:VAL:HB	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:ALA:O	1:B:181:PHE:HD2	2.04	0.40
1:B:322:LEU:O	1:B:323:VAL:CG1	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	318/334 (95%)	231 (73%)	61 (19%)	26 (8%)	<b>1</b>	<b>2</b>
1	B	315/334 (94%)	220 (70%)	68 (22%)	27 (9%)	<b>1</b>	<b>2</b>
All	All	633/668 (95%)	451 (71%)	129 (20%)	53 (8%)	<b>1</b>	<b>2</b>

All (53) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	88	ASN
1	A	100	ARG
1	A	111	ASP
1	A	144	LEU
1	A	149	ALA
1	A	167	HIS
1	A	206	ASP
1	A	248	HIS
1	A	314	GLU
1	B	51	ASN
1	B	88	ASN
1	B	122	GLU
1	B	156	ASP
1	B	184	HIS
1	B	206	ASP

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Mol	Chain	Res	Type
1	B	236	GLU
1	B	248	HIS
1	B	249	HIS
1	B	285	ARG
1	A	17	VAL
1	A	143	GLY
1	A	166	VAL
1	A	184	HIS
1	A	208	ALA
1	A	261	GLY
1	B	64	GLU
1	B	113	VAL
1	B	247	LYS
1	B	267	ASP
1	B	298	PHE
1	B	306	ARG
1	B	313	GLY
1	B	314	GLU
1	A	105	ASP
1	A	148	HIS
1	B	49	ALA
1	B	70	ALA
1	A	146	HIS
1	A	211	TYR
1	A	82	GLN
1	A	89	ASN
1	A	299	ASP
1	B	17	VAL
1	B	99	ASP
1	B	190	ASP
1	A	70	ALA
1	B	94	ASP
1	A	93	GLY
1	B	96	VAL
1	A	173	GLY
1	B	128	GLY
1	B	262	PRO

### 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	264/277 (95%)	243 (92%)	21 (8%)	14	38
1	B	264/277 (95%)	246 (93%)	18 (7%)	18	47
All	All	528/554 (95%)	489 (93%)	39 (7%)	16	42

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	22	LYS
1	A	24	VAL
1	A	29	VAL
1	A	35	PHE
1	A	53	LEU
1	A	86	LEU
1	A	133	LEU
1	A	142	GLN
1	A	176	THR
1	A	188	VAL
1	A	231	SER
1	A	243	ASN
1	A	267	ASP
1	A	270	LEU
1	A	276	CYS
1	A	292	ARG
1	A	299	ASP
1	A	302	THR
1	A	306	ARG
1	A	309	LYS
1	B	3	LYS
1	B	22	LYS
1	B	35	PHE
1	B	56	GLN
1	B	119	SER
1	B	133	LEU
1	B	188	VAL
1	B	205	CYS
1	B	206	ASP
1	B	223	ASP

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Mol	Chain	Res	Type
1	B	243	ASN
1	B	251	LEU
1	B	267	ASP
1	B	292	ARG
1	B	299	ASP
1	B	306	ARG
1	B	309	LYS
1	B	326	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	56	GLN
1	A	77	ASN
1	A	82	GLN
1	A	95	ASN
1	A	116	GLN
1	A	134	ASN
1	A	142	GLN
1	A	213	HIS
1	A	237	HIS
1	A	243	ASN
1	A	249	HIS
1	A	255	HIS
1	A	286	GLN
1	A	301	ASN
1	B	16	GLN
1	B	28	GLN
1	B	56	GLN
1	B	77	ASN
1	B	88	ASN
1	B	116	GLN
1	B	134	ASN
1	B	170	GLN
1	B	213	HIS
1	B	237	HIS
1	B	243	ASN
1	B	249	HIS
1	B	255	HIS
1	B	274	HIS
1	B	286	GLN

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Mol	Chain	Res	Type
1	B	301	ASN
1	B	304	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 6 ligands modelled in this entry, 6 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

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	320/334 (95%)	0.12	15 (4%) 32 22	17, 34, 58, 67	0
1	B	319/334 (95%)	-0.02	7 (2%) 62 52	18, 34, 58, 68	0
All	All	639/668 (95%)	0.05	22 (3%) 46 34	17, 34, 58, 68	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	ASP	4.4
1	A	149	ALA	4.2
1	A	298	PHE	3.8
1	A	148	HIS	3.8
1	A	19	GLY	3.5
1	B	300	ASP	3.3
1	A	122	GLU	3.3
1	A	145	THR	3.1
1	A	18	ASP	2.7
1	B	93	GLY	2.6
1	A	142	GLN	2.5
1	B	298	PHE	2.4
1	B	185	ARG	2.4
1	B	16	GLN	2.3
1	A	21	LYS	2.3
1	B	11	ARG	2.3
1	A	146	HIS	2.3
1	A	322	LEU	2.1
1	A	93	GLY	2.1
1	B	261	GLY	2.1
1	A	50	ASP	2.0
1	A	141	ILE	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 carbohydrates 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. 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	CS	B	403	1/1	0.98	0.15	-1.60	27,27,27,27	1
2	CS	A	400	1/1	0.92	0.12	-1.93	27,27,27,27	1
2	CS	B	405	1/1	0.47	0.31	-	27,27,27,27	1
2	CS	B	404	1/1	0.86	0.29	-	27,27,27,27	1
2	CS	A	401	1/1	0.84	0.12	-	27,27,27,27	1
2	CS	A	402	1/1	0.25	0.55	-	27,27,27,27	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.