



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:36 AM GMT

PDB ID : 2B0U  
Title : The Structure of the Follistatin:Activin Complex  
Authors : Thompson, T.B.; Lerch, T.F.; Cook, R.W.; Woodruff, T.K.; Jardetzky, T.S.  
Deposited on : 2005-09-14  
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  
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

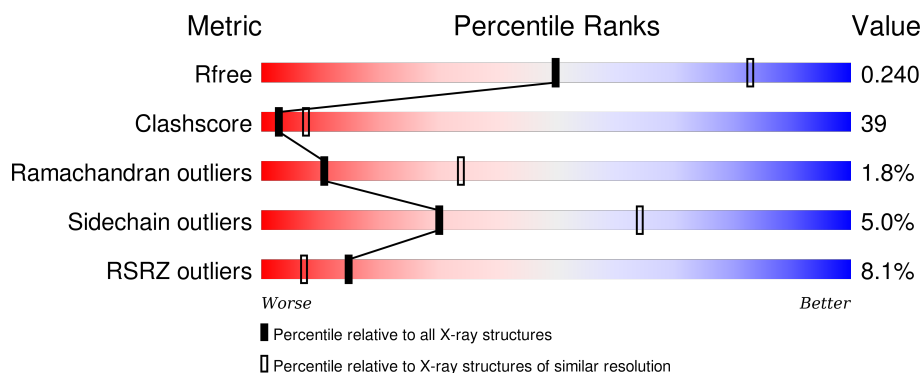
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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	116	<div> <div>3%</div> <div>55%</div> <div>38%</div> <div>7%</div> </div>
1	B	116	<div> <div>3%</div> <div>53%</div> <div>44%</div> <div>.</div> </div>
2	C	288	<div> <div>9%</div> <div>55%</div> <div>42%</div> <div>.</div> </div>
2	D	288	<div> <div>11%</div> <div>51%</div> <div>42%</div> <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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MLI	A	402	-	-	-	X
4	MLI	A	403	-	-	-	X
4	MLI	C	400	-	-	-	X
5	MPD	B	404	-	-	X	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6045 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 Inhibin beta A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	0	0
			883	553	150	167	13			
1	B	116	Total	C	N	O	S	0	0	0
			882	554	151	164	13			

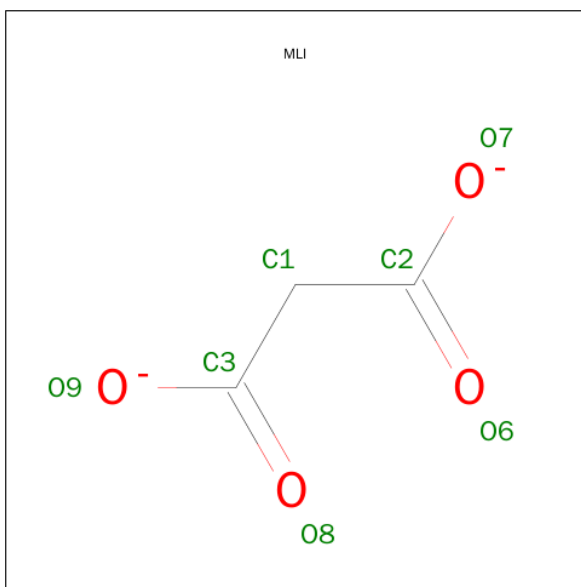
- Molecule 2 is a protein called Follistatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	287	Total	C	N	O	S	0	0	0
			2069	1251	365	414	39			
2	D	280	Total	C	N	O	S	0	0	0
			2029	1230	354	406	39			

- Molecule 3 is IRIDIUM (III) ION (three-letter code: IR3) (formula: Ir).

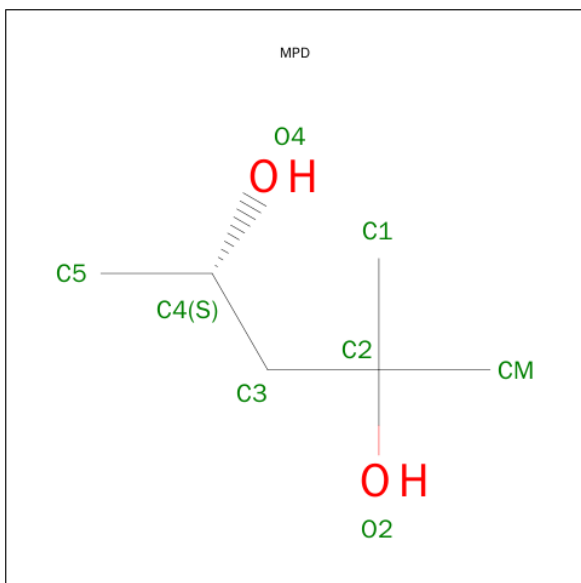
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ir	0	0
			1	1		
3	A	1	Total	Ir	0	0
			1	1		
3	D	1	Total	Ir	0	0
			1	1		
3	C	2	Total	Ir	0	0
			2	2		

- Molecule 4 is MALONATE ION (three-letter code: MLI) (formula: C<sub>3</sub>H<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			7	3	4		
4	A	1	Total	C	O	0	0
			7	3	4		
4	A	1	Total	C	O	0	0
			7	3	4		
4	A	1	Total	C	O	0	0
			7	3	4		

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:  $C_6H_{14}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			8	6	2		

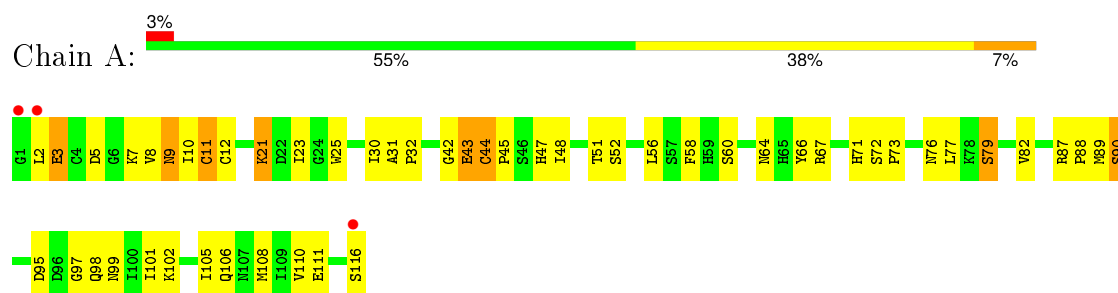
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	25	Total	O	0	0
			25	25		
6	B	31	Total	O	0	0
			31	31		
6	C	48	Total	O	0	0
			48	48		
6	D	37	Total	O	0	0
			37	37		

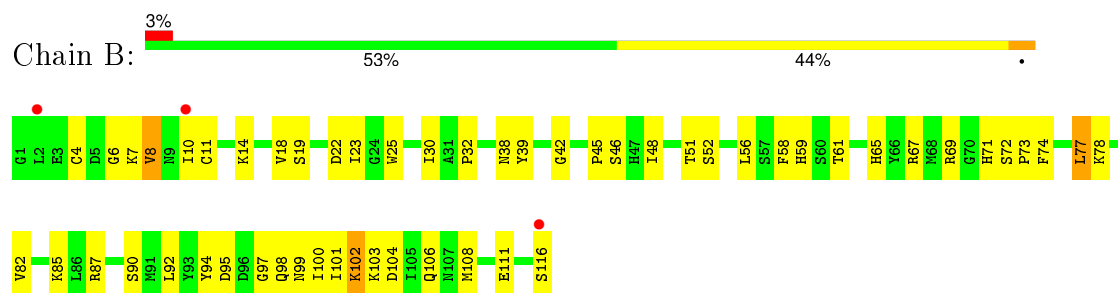
### 3 Residue-property plots [i](#)

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.

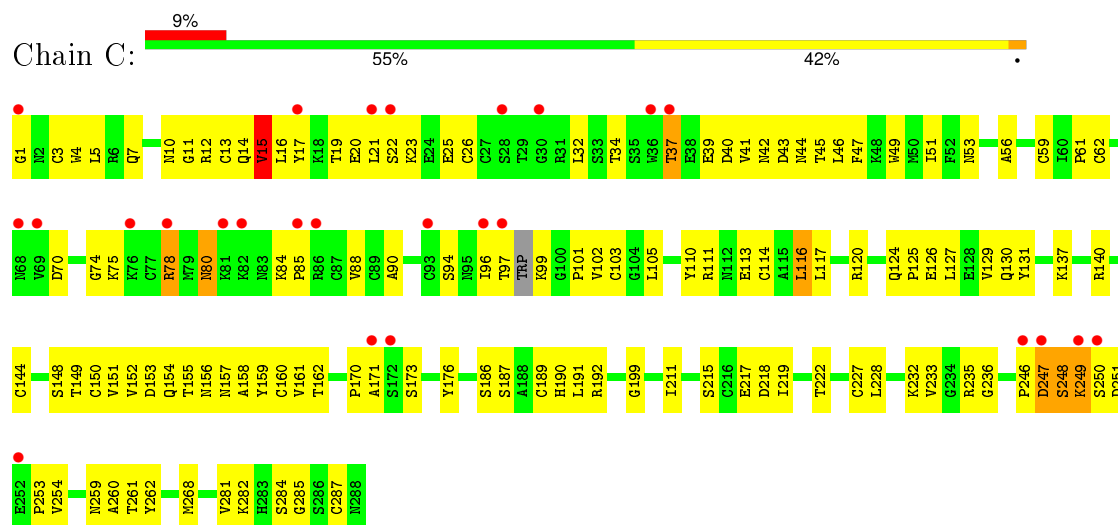
#### • Molecule 1: Inhibin beta A chain



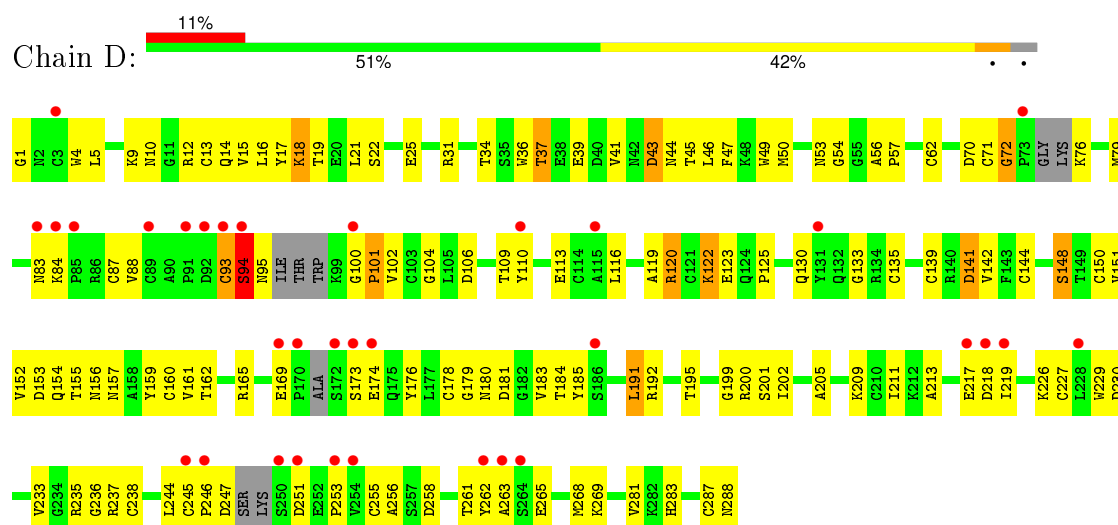
#### • Molecule 1: Inhibin beta A chain



#### • Molecule 2: Follistatin



#### • Molecule 2: Follistatin





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.25Å 120.82Å 70.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.35 – 2.80 29.64 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.35-2.80) 99.0 (29.64-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.88 (at 2.80Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.265 , 0.297 0.269 , 0.240	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	68.6	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 57.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 23595 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6045	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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, 0.375 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: MPD, MLI, IR3

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	0.58	0/905	0.86	0/1223
1	B	0.59	0/905	0.92	2/1223 (0.2%)
2	C	0.57	0/2100	0.91	1/2839 (0.0%)
2	D	0.53	0/2058	0.92	3/2781 (0.1%)
All	All	0.56	0/5968	0.91	6/8066 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	93	CYS	CA-CB-SG	7.42	127.36	114.00
2	D	72	GLY	N-CA-C	-6.98	95.64	113.10
2	D	101	PRO	N-CA-C	-6.21	95.95	112.10
1	B	78	LYS	N-CA-C	5.90	126.93	111.00
1	B	8	VAL	N-CA-C	-5.64	95.78	111.00
2	C	80	ASN	N-CA-C	5.36	125.48	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	883	0	816	72	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	882	0	815	77	0
2	C	2069	0	1895	181	1
2	D	2029	0	1850	159	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
4	A	21	0	6	0	0
4	C	7	0	2	0	0
5	B	8	0	14	9	0
6	A	25	0	0	3	0
6	B	31	0	0	3	0
6	C	48	0	0	7	0
6	D	37	0	0	3	0
All	All	6045	0	5398	443	1

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 39.

All (443) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:261:THR:HG22	2:C:287:CYS:N	1.56	1.18
2:D:120:ARG:HG2	2:D:120:ARG:HH11	1.10	1.15
2:C:261:THR:CG2	2:C:287:CYS:H	1.60	1.14
2:C:21:LEU:HD11	2:C:25:GLU:HB2	1.24	1.12
1:B:56:LEU:HD13	1:B:116:SER:HB3	1.28	1.09
2:C:19:THR:HG22	2:C:20:GLU:HG3	1.37	1.06
2:C:78:ARG:HG3	2:C:78:ARG:HH11	1.21	1.06
2:C:21:LEU:HD21	2:C:26:CYS:HB2	1.38	1.03
1:A:56:LEU:HD13	1:A:116:SER:CB	1.94	0.98
1:A:8:VAL:HG12	1:A:10:ILE:HG13	1.46	0.98
1:B:30:ILE:HD11	1:B:94:TYR:CE1	1.98	0.97
2:D:5:LEU:HD22	2:D:17:TYR:OH	1.63	0.97
2:C:21:LEU:HD12	2:C:22:SER:H	1.25	0.97
2:C:173:SER:HA	2:C:176:TYR:CD1	1.99	0.97
2:D:261:THR:HG22	2:D:287:CYS:H	1.29	0.96
1:A:10:ILE:HG23	1:A:48:ILE:HD13	1.45	0.95
2:C:21:LEU:CD1	2:C:25:GLU:HB2	1.95	0.95
2:D:144:CYS:HB3	2:D:148:SER:OG	1.66	0.95
2:D:4:TRP:HE1	2:D:37:THR:HB	1.33	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:ASP:HB3	1:B:101:ILE:HD12	1.51	0.92
2:D:142:VAL:HG11	2:D:150:CYS:HB2	1.52	0.91
1:A:2:LEU:HD13	2:D:5:LEU:HD23	1.52	0.89
2:C:19:THR:HG22	2:C:20:GLU:CG	2.03	0.89
1:A:98:GLN:HE22	2:D:192:ARG:HE	1.19	0.89
1:B:98:GLN:HE22	2:C:192:ARG:HE	0.90	0.89
1:B:98:GLN:HE22	2:C:192:ARG:NE	1.71	0.89
1:A:45:PRO:HD2	1:A:48:ILE:HD12	1.56	0.88
2:C:261:THR:HG22	2:C:287:CYS:H	0.73	0.88
2:C:78:ARG:CG	2:C:78:ARG:HH11	1.87	0.88
5:B:404:MPD:H12	2:C:105:LEU:HG	1.54	0.87
2:D:109:THR:HG22	2:D:135:CYS:H	1.39	0.86
2:D:120:ARG:HG2	2:D:120:ARG:NH1	1.85	0.85
2:D:211:ILE:HD11	2:D:219:ILE:HG12	1.56	0.85
2:C:74:GLY:O	2:C:90:ALA:N	2.07	0.85
2:D:120:ARG:CG	2:D:120:ARG:HH11	1.87	0.85
2:C:4:TRP:HE1	2:C:37:THR:HB	1.42	0.84
1:A:2:LEU:CD1	2:D:5:LEU:CD2	2.55	0.84
1:A:8:VAL:O	1:A:9:ASN:HB3	1.76	0.83
2:D:5:LEU:HB2	2:D:17:TYR:CE1	2.13	0.83
2:C:41:VAL:HG11	2:C:45:THR:HG21	1.61	0.83
2:C:261:THR:HG23	2:C:287:CYS:HB2	1.60	0.82
1:B:10:ILE:HG23	1:B:48:ILE:HD13	1.61	0.82
1:A:51:THR:HG21	2:D:16:LEU:HB2	1.61	0.82
2:D:5:LEU:HD13	2:D:17:TYR:OH	1.80	0.82
2:C:173:SER:HA	2:C:176:TYR:HD1	1.41	0.82
1:B:10:ILE:CG2	1:B:48:ILE:HD13	2.11	0.81
1:B:98:GLN:NE2	2:C:192:ARG:HE	1.74	0.81
1:A:98:GLN:NE2	2:D:192:ARG:HE	1.79	0.80
2:C:268:MET:CE	2:C:281:VAL:HG22	2.12	0.80
2:D:161:VAL:HG12	2:D:162:THR:N	1.97	0.80
2:C:215:SER:OG	2:C:217:GLU:HG3	1.82	0.79
1:B:51:THR:HG21	2:C:16:LEU:H	1.47	0.79
2:C:261:THR:CG2	2:C:287:CYS:HB2	2.12	0.79
1:B:65:HIS:HB3	1:B:69:ARG:NH1	1.98	0.78
1:A:66:TYR:HE1	1:B:18:VAL:HG13	1.48	0.78
1:B:56:LEU:HD13	1:B:116:SER:CB	2.10	0.78
1:A:56:LEU:HD13	1:A:116:SER:HB3	1.64	0.78
1:A:10:ILE:CG2	1:A:48:ILE:HD13	2.14	0.77
1:B:45:PRO:HG2	1:B:48:ILE:HD12	1.65	0.77
2:D:268:MET:CE	2:D:281:VAL:HG22	2.15	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:253:PRO:HA	2:D:263:ALA:O	1.83	0.77
2:C:41:VAL:HG11	2:C:45:THR:CG2	2.16	0.76
2:C:232:LYS:HE3	6:C:538:HOH:O	1.85	0.76
2:D:12:ARG:NH1	2:D:53:ASN:HB3	2.01	0.75
2:C:21:LEU:CD1	2:C:22:SER:H	1.98	0.75
1:A:108:MET:HG2	1:B:58:PHE:CE2	2.21	0.74
2:C:173:SER:HA	2:C:176:TYR:CE1	2.21	0.74
2:C:232:LYS:CE	6:C:538:HOH:O	2.34	0.74
2:D:41:VAL:HG21	2:D:49:TRP:CZ3	2.22	0.74
2:D:261:THR:HG22	2:D:287:CYS:N	2.02	0.73
2:C:41:VAL:HG11	2:C:45:THR:CB	2.18	0.73
1:A:2:LEU:CD1	2:D:5:LEU:HD23	2.18	0.73
2:C:268:MET:HE3	2:C:281:VAL:HG22	1.68	0.73
1:A:2:LEU:HD13	2:D:5:LEU:CD2	2.19	0.73
2:D:4:TRP:CG	2:D:13:CYS:HB3	2.24	0.73
2:C:14:GLN:C	2:C:15:VAL:HG23	2.09	0.72
2:D:31:ARG:HH11	2:D:31:ARG:HG3	1.55	0.72
1:A:45:PRO:CD	1:A:48:ILE:HD12	2.19	0.72
2:D:251:ASP:O	2:D:253:PRO:HD3	1.89	0.72
2:D:5:LEU:HD11	2:D:31:ARG:HE	1.56	0.71
1:B:51:THR:CG2	2:C:16:LEU:H	2.04	0.70
2:C:186:SER:HB2	2:C:190:HIS:CD2	2.26	0.70
1:B:11:CYS:HA	1:B:42:GLY:HA3	1.73	0.70
5:B:404:MPD:C1	2:C:105:LEU:HG	2.20	0.70
2:C:173:SER:CA	2:C:176:TYR:HD1	2.04	0.70
1:A:51:THR:HG21	2:D:16:LEU:CB	2.21	0.70
1:A:48:ILE:O	1:A:52:SER:HB2	1.91	0.69
2:C:251:ASP:O	2:C:253:PRO:HD3	1.93	0.69
1:A:47:HIS:O	1:A:51:THR:HB	1.93	0.69
1:B:23:ILE:HG22	1:B:25:TRP:CD1	2.28	0.69
1:A:8:VAL:CG1	1:A:10:ILE:HG13	2.23	0.69
2:D:184:THR:HG23	2:D:209:LYS:HA	1.74	0.69
2:C:41:VAL:CG1	2:C:45:THR:HB	2.23	0.69
2:D:261:THR:CG2	2:D:287:CYS:HB2	2.22	0.68
1:B:14:LYS:O	1:B:38:ASN:HB3	1.94	0.67
1:A:64:ASN:OD1	1:A:67:ARG:NH1	2.28	0.67
2:D:5:LEU:CD2	2:D:17:TYR:OH	2.41	0.66
2:C:21:LEU:CD1	2:C:25:GLU:CB	2.73	0.66
2:C:80:ASN:HB2	2:C:84:LYS:H	1.60	0.66
2:C:102:VAL:HG11	2:C:129:VAL:HG22	1.75	0.66
2:C:3:CYS:HB2	2:C:21:LEU:HD23	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:5:LEU:HB2	2:D:17:TYR:HE1	1.59	0.66
2:D:5:LEU:O	2:D:5:LEU:HG	1.94	0.66
2:C:249:LYS:C	2:C:251:ASP:H	1.96	0.66
2:C:187:SER:OG	2:C:190:HIS:CD2	2.49	0.65
1:A:95:ASP:HB3	1:A:101:ILE:HD12	1.78	0.65
1:B:30:ILE:HD11	1:B:94:TYR:CZ	2.31	0.65
2:D:195:THR:OG1	2:D:202:ILE:HD12	1.96	0.65
2:D:109:THR:CG2	2:D:135:CYS:H	2.07	0.65
2:C:41:VAL:HG12	2:C:42:ASN:N	2.10	0.65
1:A:21:LYS:HB3	1:A:21:LYS:NZ	2.12	0.65
2:C:21:LEU:HD12	2:C:22:SER:N	2.07	0.65
2:C:102:VAL:HG12	2:C:103:CYS:N	2.12	0.64
2:D:261:THR:HG23	2:D:287:CYS:HB2	1.79	0.64
1:A:58:PHE:CD2	1:B:108:MET:HG2	2.32	0.64
1:A:10:ILE:HG23	1:A:48:ILE:CD1	2.26	0.64
2:D:39:GLU:HG2	2:D:41:VAL:HG13	1.79	0.64
2:D:161:VAL:CG1	2:D:162:THR:N	2.61	0.64
2:D:144:CYS:HB3	2:D:148:SER:CB	2.28	0.63
1:B:10:ILE:HG22	1:B:11:CYS:H	1.62	0.63
1:B:95:ASP:OD2	1:B:101:ILE:HD11	1.98	0.63
1:B:51:THR:HG21	2:C:16:LEU:HB3	1.80	0.63
2:C:246:PRO:HG3	2:D:9:LYS:O	1.99	0.63
2:D:4:TRP:NE1	2:D:37:THR:HB	2.10	0.63
2:D:130:GLN:NE2	2:D:160:CYS:H	1.96	0.63
1:A:72:SER:HB2	1:A:73:PRO:HD3	1.80	0.63
2:C:211:ILE:HD11	2:C:218:ASP:O	1.97	0.63
2:C:247:ASP:O	2:C:248:SER:OG	2.14	0.62
1:A:45:PRO:HG2	1:A:48:ILE:HD12	1.80	0.62
2:D:141:ASP:C	2:D:142:VAL:CG2	2.68	0.62
2:D:180:ASN:O	2:D:235:ARG:NH1	2.32	0.62
2:D:71:CYS:SG	2:D:76:LYS:O	2.56	0.62
1:B:65:HIS:HB3	1:B:69:ARG:HH12	1.63	0.62
2:C:11:GLY:O	2:C:59:CYS:HB3	1.98	0.62
2:C:130:GLN:NE2	2:C:160:CYS:H	1.98	0.62
1:A:23:ILE:HG22	1:A:25:TRP:CD1	2.34	0.62
1:A:105:ILE:HD12	2:C:51:ILE:HG23	1.81	0.62
2:D:269:LYS:NZ	6:D:537:HOH:O	2.32	0.62
2:C:219:ILE:CD1	2:C:236:GLY:HA3	2.30	0.62
1:B:45:PRO:HG2	1:B:48:ILE:CD1	2.30	0.61
1:A:66:TYR:CE1	1:B:18:VAL:HG13	2.34	0.61
2:C:154:GLN:HG2	2:C:155:THR:HG23	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:256:ALA:HB3	2:D:258:ASP:OD1	2.00	0.61
2:C:49:TRP:CD1	2:C:53:ASN:ND2	2.69	0.61
1:B:14:LYS:HD2	1:B:39:TYR:CE1	2.35	0.61
2:C:21:LEU:CG	2:C:22:SER:N	2.63	0.61
1:B:8:VAL:HG12	1:B:10:ILE:CD1	2.30	0.61
1:A:56:LEU:HD13	1:A:116:SER:OG	2.01	0.61
2:D:130:GLN:HE22	2:D:159:TYR:HA	1.66	0.60
2:C:5:LEU:CB	2:C:17:TYR:HE1	2.14	0.60
1:A:8:VAL:O	1:A:9:ASN:CB	2.48	0.60
2:C:5:LEU:CB	2:C:15:VAL:HB	2.32	0.60
1:A:31:ALA:HA	1:A:32:PRO:O	2.02	0.60
2:C:21:LEU:HD11	2:C:25:GLU:CB	2.15	0.60
2:C:173:SER:CA	2:C:176:TYR:CD1	2.79	0.60
1:B:45:PRO:CG	1:B:48:ILE:HD12	2.32	0.60
2:C:7:GLN:HE22	2:C:61:PRO:HG3	1.66	0.60
2:C:21:LEU:HG	2:C:22:SER:N	2.17	0.60
1:B:10:ILE:HG22	1:B:11:CYS:N	2.17	0.60
2:D:10:ASN:HD21	2:D:12:ARG:HD3	1.65	0.60
2:D:268:MET:HE1	2:D:281:VAL:HG22	1.83	0.59
2:C:41:VAL:HG11	2:C:45:THR:HB	1.82	0.59
2:D:123:GLU:O	2:D:125:PRO:HD3	2.03	0.59
1:B:51:THR:HG23	2:C:4:TRP:HZ3	1.66	0.59
1:A:45:PRO:CG	1:A:48:ILE:HD12	2.33	0.59
2:D:44:ASN:OD1	2:D:47:PHE:HB3	2.02	0.59
2:D:169:GLU:HB2	6:D:527:HOH:O	2.02	0.59
1:A:95:ASP:OD2	1:A:101:ILE:HD11	2.03	0.58
2:C:46:LEU:O	2:C:49:TRP:HB2	2.04	0.58
2:C:259:ASN:HD21	2:C:282:LYS:NZ	2.01	0.58
2:D:76:LYS:O	2:D:87:CYS:HA	2.03	0.58
1:A:87:ARG:HG2	1:A:88:PRO:HD2	1.85	0.58
1:B:8:VAL:HG11	2:C:17:TYR:HE2	1.68	0.58
2:C:102:VAL:HG11	2:C:129:VAL:CG2	2.32	0.58
1:B:71:HIS:CE1	6:B:531:HOH:O	2.55	0.58
2:D:200:ARG:HG2	2:D:201:SER:N	2.18	0.58
2:C:21:LEU:HD12	2:C:25:GLU:OE2	2.04	0.58
2:D:79:MET:CE	2:D:83:ASN:OD1	2.52	0.58
1:B:85:LYS:O	1:B:111:GLU:HB3	2.02	0.58
2:C:21:LEU:HD12	2:C:25:GLU:CD	2.25	0.57
2:D:18:LYS:HG3	2:D:19:THR:H	1.67	0.57
1:B:67:ARG:HH21	1:B:77:LEU:HB3	1.68	0.57
2:C:41:VAL:CG1	2:C:45:THR:CB	2.82	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:4:TRP:HH2	2:D:46:LEU:HD22	1.69	0.57
2:D:141:ASP:O	2:D:142:VAL:HG22	2.03	0.57
1:B:90:SER:O	5:B:404:MPD:HM1	2.04	0.57
2:C:268:MET:HE1	2:C:281:VAL:HG22	1.86	0.57
2:C:246:PRO:HG3	2:D:9:LYS:C	2.25	0.57
2:C:5:LEU:CA	2:C:17:TYR:HE1	2.18	0.57
1:A:95:ASP:OD1	1:A:97:GLY:N	2.38	0.57
2:C:268:MET:HE1	2:C:281:VAL:HA	1.87	0.57
2:C:78:ARG:HG3	2:C:78:ARG:NH1	2.01	0.56
2:D:31:ARG:NH1	2:D:31:ARG:HG3	2.20	0.56
2:D:139:CYS:O	2:D:142:VAL:CG2	2.53	0.56
2:D:174:GLU:O	2:D:174:GLU:HG2	2.05	0.56
2:C:144:CYS:HB3	2:C:148:SER:OG	2.06	0.56
1:B:51:THR:CG2	2:C:4:TRP:HZ3	2.19	0.56
2:C:102:VAL:CG1	2:C:129:VAL:HG13	2.35	0.56
1:A:98:GLN:HE22	2:D:192:ARG:NE	1.98	0.56
1:B:90:SER:C	5:B:404:MPD:HM1	2.26	0.56
2:D:161:VAL:HG12	2:D:162:THR:H	1.69	0.56
2:D:155:THR:HG22	2:D:155:THR:O	2.06	0.56
2:D:130:GLN:HE21	2:D:160:CYS:H	1.52	0.56
2:D:5:LEU:CD1	2:D:17:TYR:OH	2.53	0.55
1:B:103:LYS:HG2	1:B:104:ASP:N	2.20	0.55
1:B:61:THR:HG21	2:C:47:PHE:CE1	2.41	0.55
2:C:21:LEU:CD1	2:C:22:SER:N	2.67	0.55
2:D:100:GLY:O	2:D:102:VAL:HG13	2.06	0.55
2:D:21:LEU:HD22	2:D:25:GLU:OE1	2.07	0.55
1:A:9:ASN:CG	1:A:9:ASN:O	2.45	0.55
2:C:170:PRO:O	2:C:171:ALA:HB3	2.07	0.55
2:D:213:ALA:HB1	2:D:218:ASP:HB2	1.89	0.55
2:C:12:ARG:HG2	2:C:56:ALA:O	2.06	0.55
1:A:82:VAL:HG21	1:B:82:VAL:HG21	1.89	0.55
2:C:151:VAL:HG22	2:C:199:GLY:HA2	1.89	0.55
2:C:161:VAL:HG12	2:C:162:THR:N	2.22	0.54
1:B:102:LYS:C	1:B:102:LYS:HD2	2.27	0.54
2:C:41:VAL:HG12	2:C:45:THR:HB	1.89	0.54
2:D:161:VAL:CG1	2:D:162:THR:H	2.19	0.54
2:C:97:THR:O	2:C:99:LYS:HB2	2.07	0.54
1:B:10:ILE:HG23	1:B:48:ILE:CD1	2.34	0.54
1:A:56:LEU:CD1	1:A:116:SER:HB3	2.34	0.54
1:A:2:LEU:HD11	2:D:5:LEU:CD2	2.36	0.54
1:B:95:ASP:HB3	1:B:101:ILE:CD1	2.30	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:130:GLN:NE2	2:D:159:TYR:HA	2.23	0.54
2:D:153:ASP:OD2	2:D:157:ASN:HB2	2.07	0.54
1:A:47:HIS:O	1:A:51:THR:CB	2.55	0.53
2:D:10:ASN:HD21	2:D:12:ARG:CD	2.21	0.53
2:D:34:THR:O	2:D:62:CYS:HB2	2.08	0.53
2:C:117:LEU:CD2	2:C:120:ARG:HH21	2.20	0.53
2:C:32:LEU:HD11	2:C:85:PRO:HB2	1.90	0.53
2:C:110:TYR:CD1	2:C:116:LEU:HD12	2.43	0.53
2:C:78:ARG:CG	2:C:78:ARG:NH1	2.55	0.52
1:B:102:LYS:HG3	2:C:159:TYR:OH	2.08	0.52
2:D:83:ASN:CG	2:D:83:ASN:O	2.44	0.52
2:C:116:LEU:HD11	2:C:127:LEU:HG	1.91	0.52
2:C:232:LYS:HE2	6:C:538:HOH:O	2.04	0.52
2:C:260:ALA:O	2:C:262:TYR:CD1	2.62	0.52
2:D:219:ILE:CD1	2:D:236:GLY:HA3	2.39	0.52
2:D:5:LEU:HD11	2:D:31:ARG:NE	2.21	0.52
2:D:151:VAL:HG22	2:D:199:GLY:HA2	1.90	0.52
1:B:90:SER:O	5:B:404:MPD:CM	2.57	0.52
1:A:58:PHE:CD1	2:D:47:PHE:HZ	2.28	0.52
2:C:4:TRP:CG	2:C:13:CYS:HB3	2.44	0.52
2:C:4:TRP:O	2:C:34:THR:HG23	2.09	0.52
2:C:260:ALA:HB3	2:C:262:TYR:CE1	2.44	0.52
2:C:187:SER:OG	2:C:189:CYS:HB3	2.10	0.52
2:C:12:ARG:NH1	2:C:53:ASN:O	2.43	0.52
2:D:4:TRP:CD2	2:D:13:CYS:HB3	2.45	0.51
2:C:39:GLU:HG2	2:C:41:VAL:CG2	2.40	0.51
2:C:1:GLY:HA2	2:C:40:ASP:OD1	2.10	0.51
1:B:30:ILE:HB	1:B:92:LEU:O	2.10	0.51
2:C:5:LEU:CB	2:C:17:TYR:CE1	2.94	0.51
1:A:31:ALA:HA	1:A:32:PRO:C	2.31	0.51
2:C:249:LYS:C	2:C:251:ASP:N	2.64	0.51
2:C:34:THR:O	2:C:62:CYS:HB2	2.10	0.50
2:C:219:ILE:HD11	2:C:236:GLY:HA3	1.92	0.50
2:C:78:ARG:HH11	2:C:88:VAL:HG21	1.77	0.50
1:B:32:PRO:HD3	5:B:404:MPD:HM2	1.92	0.50
2:D:261:THR:HG22	2:D:287:CYS:CB	2.42	0.50
1:A:51:THR:HG22	2:D:16:LEU:H	1.77	0.50
2:D:79:MET:HE3	2:D:83:ASN:OD1	2.11	0.50
2:C:44:ASN:O	2:C:47:PHE:HB3	2.12	0.50
1:A:105:ILE:CD1	2:C:51:ILE:HG23	2.42	0.50
1:A:45:PRO:HG2	1:A:48:ILE:CD1	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:154:GLN:HG2	2:C:155:THR:N	2.27	0.49
2:D:154:GLN:HG2	2:D:155:THR:N	2.26	0.49
1:B:48:ILE:O	1:B:52:SER:HB2	2.12	0.49
1:B:14:LYS:HD2	1:B:39:TYR:CZ	2.47	0.49
2:D:181:ASP:C	2:D:235:ARG:NH1	2.66	0.49
2:D:1:GLY:HA3	2:D:36:TRP:NE1	2.27	0.49
5:B:404:MPD:C5	2:C:126:GLU:HB2	2.42	0.49
1:B:100:ILE:HD13	2:C:151:VAL:HG13	1.94	0.49
2:C:102:VAL:CG1	2:C:103:CYS:N	2.76	0.49
1:A:76:ASN:ND2	6:A:512:HOH:O	2.45	0.49
2:C:211:ILE:CD1	2:C:218:ASP:O	2.60	0.49
2:C:152:VAL:HA	2:C:157:ASN:O	2.12	0.49
2:C:187:SER:HG	2:C:189:CYS:HB3	1.77	0.49
2:D:180:ASN:C	2:D:235:ARG:HH12	2.15	0.48
2:C:149:THR:HG22	2:C:150:CYS:N	2.28	0.48
2:C:170:PRO:HG3	2:C:186:SER:HB3	1.95	0.48
2:D:265:GLU:O	2:D:268:MET:HB3	2.13	0.48
2:C:249:LYS:HB3	6:C:540:HOH:O	2.13	0.48
1:A:30:ILE:HG23	2:D:161:VAL:HG11	1.95	0.48
1:B:30:ILE:HG23	2:C:161:VAL:HG11	1.96	0.48
2:C:39:GLU:HG2	2:C:41:VAL:HG23	1.96	0.48
2:C:187:SER:HG	2:C:190:HIS:CD2	2.31	0.48
1:B:19:SER:HB3	1:B:22:ASP:OD2	2.14	0.48
2:D:287:CYS:O	2:D:288:ASN:CB	2.61	0.48
1:B:98:GLN:NE2	2:C:192:ARG:HG3	2.29	0.48
2:D:261:THR:HG22	2:D:287:CYS:HB2	1.96	0.48
2:D:130:GLN:HE22	2:D:159:TYR:HD1	1.61	0.48
2:D:93:CYS:O	2:D:94:SER:C	2.51	0.48
6:A:519:HOH:O	2:D:161:VAL:HG13	2.14	0.47
2:C:41:VAL:CG1	2:C:42:ASN:N	2.76	0.47
2:D:1:GLY:HA3	2:D:36:TRP:CE2	2.49	0.47
2:C:84:LYS:HZ1	2:D:244:LEU:HD11	1.79	0.47
2:C:285:GLY:N	6:C:544:HOH:O	2.38	0.47
1:A:9:ASN:ND2	1:A:9:ASN:O	2.47	0.47
1:A:10:ILE:HG22	1:A:11:CYS:N	2.30	0.47
2:C:259:ASN:HD21	2:C:282:LYS:HZ2	1.61	0.47
2:D:178:CYS:O	2:D:205:ALA:HB3	2.14	0.47
2:C:78:ARG:HG3	2:C:88:VAL:HG21	1.97	0.47
1:A:23:ILE:HG22	1:A:25:TRP:NE1	2.30	0.47
2:C:70:ASP:N	6:C:541:HOH:O	2.39	0.47
1:A:7:LYS:CB	6:A:527:HOH:O	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:173:SER:C	2:C:176:TYR:HD1	2.19	0.46
2:C:227:CYS:O	2:C:228:LEU:HD23	2.15	0.46
2:C:261:THR:HG22	2:C:287:CYS:HB2	1.97	0.46
2:C:113:GLU:O	2:C:116:LEU:HB3	2.15	0.46
2:C:101:PRO:HA	2:C:111:ARG:O	2.15	0.46
1:B:45:PRO:CD	1:B:48:ILE:HD12	2.45	0.46
2:D:12:ARG:CG	2:D:56:ALA:O	2.64	0.46
2:D:113:GLU:O	2:D:116:LEU:HB3	2.16	0.46
2:D:151:VAL:HG12	2:D:152:VAL:N	2.31	0.46
2:C:23:LYS:O	2:C:26:CYS:HB3	2.16	0.46
2:D:44:ASN:O	2:D:47:PHE:N	2.47	0.46
1:A:88:PRO:HB3	1:A:106:GLN:HA	1.97	0.46
2:D:79:MET:HE1	2:D:83:ASN:OD1	2.14	0.46
2:D:18:LYS:HG3	2:D:19:THR:N	2.30	0.46
2:D:151:VAL:CG1	2:D:152:VAL:N	2.78	0.46
1:B:10:ILE:HG21	1:B:48:ILE:HD13	1.94	0.46
1:B:69:ARG:O	6:B:531:HOH:O	2.21	0.46
1:B:45:PRO:HD2	1:B:48:ILE:HD12	1.96	0.46
2:D:180:ASN:HB2	2:D:202:ILE:O	2.16	0.46
2:C:150:CYS:O	2:C:151:VAL:HG23	2.16	0.46
2:D:119:ALA:O	2:D:122:LYS:HB3	2.17	0.45
2:D:227:CYS:HA	2:D:237:ARG:O	2.16	0.45
2:D:70:ASP:OD1	2:D:72:GLY:N	2.49	0.45
2:C:130:GLN:HE21	2:C:160:CYS:H	1.65	0.45
2:C:249:LYS:HA	2:C:249:LYS:HD2	1.54	0.45
2:C:148:SER:HA	2:C:161:VAL:O	2.16	0.45
1:B:74:PHE:HA	1:B:77:LEU:HB2	1.99	0.45
2:D:1:GLY:HA3	2:D:36:TRP:CZ2	2.52	0.45
1:B:87:ARG:HH22	2:C:125:PRO:HD3	1.82	0.45
2:C:117:LEU:HD23	2:C:120:ARG:HH21	1.81	0.45
1:A:51:THR:HG22	1:A:51:THR:O	2.16	0.45
1:A:77:LEU:HD23	1:A:77:LEU:HA	1.76	0.45
2:D:230:ASP:OD2	2:D:233:VAL:HG22	2.16	0.45
2:C:4:TRP:HE1	2:C:37:THR:CB	2.21	0.45
1:B:72:SER:HB3	1:B:73:PRO:HD3	1.98	0.45
2:D:183:VAL:HB	2:D:185:TYR:CE1	2.52	0.45
2:D:101:PRO:O	2:D:133:GLY:N	2.42	0.45
2:C:78:ARG:CG	2:C:88:VAL:HG21	2.46	0.45
2:D:83:ASN:OD1	2:D:83:ASN:C	2.56	0.45
2:D:255:CYS:SG	2:D:283:HIS:CE1	3.10	0.45
2:C:131:TYR:HB3	2:C:160:CYS:SG	2.56	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:141:ASP:C	2:D:142:VAL:HG23	2.38	0.44
1:A:43:GLU:O	1:A:45:PRO:HD3	2.17	0.44
1:B:71:HIS:C	1:B:73:PRO:HD2	2.38	0.44
2:D:173:SER:O	2:D:176:TYR:HD1	1.99	0.44
2:D:226:LYS:O	2:D:238:CYS:HA	2.18	0.44
2:C:153:ASP:OD2	2:C:159:TYR:HE2	1.99	0.44
2:D:253:PRO:CA	2:D:263:ALA:O	2.62	0.44
2:D:49:TRP:CD1	2:D:53:ASN:ND2	2.85	0.44
2:C:75:LYS:HA	2:C:88:VAL:O	2.17	0.44
1:A:11:CYS:HA	1:A:42:GLY:HA3	1.98	0.44
5:B:404:MPD:H52	2:C:126:GLU:HB2	1.99	0.44
2:D:169:GLU:N	6:D:527:HOH:O	2.51	0.44
2:D:93:CYS:O	2:D:95:ASN:N	2.49	0.44
1:B:87:ARG:NH2	2:C:125:PRO:CD	2.81	0.44
2:C:260:ALA:N	2:C:287:CYS:SG	2.91	0.44
1:A:58:PHE:CZ	1:B:25:TRP:HZ3	2.36	0.44
2:D:195:THR:HG23	2:D:200:ARG:O	2.18	0.44
2:D:230:ASP:HB3	2:D:233:VAL:HG22	1.99	0.44
2:C:151:VAL:O	2:C:158:ALA:HA	2.17	0.44
2:C:3:CYS:HB2	2:C:21:LEU:CD2	2.47	0.44
1:B:97:GLY:O	1:B:98:GLN:HB2	2.18	0.44
1:A:110:VAL:HG23	1:B:59:HIS:CD2	2.53	0.43
2:D:76:LYS:CB	2:D:88:VAL:O	2.66	0.43
2:C:284:SER:HA	6:C:544:HOH:O	2.17	0.43
2:C:41:VAL:HG12	2:C:42:ASN:H	1.79	0.43
2:D:184:THR:CG2	2:D:209:LYS:HA	2.47	0.43
2:D:50:MET:O	2:D:54:GLY:HA2	2.18	0.43
2:D:139:CYS:O	2:D:142:VAL:HG23	2.19	0.43
2:D:44:ASN:O	2:D:45:THR:C	2.57	0.43
2:C:233:VAL:CG1	2:C:233:VAL:O	2.65	0.43
1:A:56:LEU:O	1:A:60:SER:HB2	2.19	0.43
2:D:104:GLY:C	2:D:106:ASP:H	2.22	0.43
2:D:181:ASP:CA	2:D:235:ARG:HH12	2.32	0.43
2:D:245:CYS:HA	2:D:246:PRO:HD3	1.84	0.43
2:C:19:THR:HG22	2:C:20:GLU:HG2	1.95	0.42
2:D:191:LEU:HD21	2:D:202:ILE:HG21	2.01	0.42
2:C:32:LEU:CD1	2:C:85:PRO:HB2	2.49	0.42
2:C:14:GLN:C	2:C:15:VAL:CG2	2.77	0.42
2:D:179:GLY:C	2:D:181:ASP:H	2.23	0.42
1:A:10:ILE:O	1:A:12:CYS:N	2.52	0.42
2:C:161:VAL:CG1	2:C:162:THR:N	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:217:GLU:HG3	2:D:217:GLU:H	1.47	0.42
2:C:137:LYS:HD2	2:C:156:ASN:O	2.20	0.42
2:C:19:THR:HG22	2:C:20:GLU:N	2.35	0.42
2:C:10:ASN:O	2:D:244:LEU:N	2.52	0.42
2:D:195:THR:HA	2:D:202:ILE:CD1	2.50	0.42
1:B:100:ILE:HD13	2:C:151:VAL:CG1	2.49	0.42
2:D:226:LYS:HB3	2:D:226:LYS:HE2	1.85	0.42
1:B:99:ASN:H	1:B:99:ASN:HD22	1.68	0.42
2:C:130:GLN:NE2	2:C:159:TYR:HA	2.35	0.42
1:A:95:ASP:OD2	1:A:99:ASN:HB2	2.19	0.42
2:D:191:LEU:HD21	2:D:202:ILE:CG2	2.49	0.42
2:D:258:ASP:OD2	2:D:262:TYR:OH	2.34	0.42
2:C:149:THR:CG2	2:C:150:CYS:N	2.83	0.42
1:B:72:SER:O	1:B:73:PRO:C	2.57	0.42
2:D:41:VAL:HG21	2:D:49:TRP:CH2	2.55	0.42
2:C:110:TYR:CE1	2:C:116:LEU:HD12	2.54	0.42
1:B:8:VAL:HG12	1:B:10:ILE:HD12	2.02	0.42
2:C:43:ASP:O	2:C:46:LEU:HB2	2.20	0.42
5:B:404:MPD:H12	2:C:105:LEU:CG	2.37	0.42
2:D:161:VAL:HG12	2:D:162:THR:O	2.20	0.42
2:D:10:ASN:ND2	2:D:12:ARG:CD	2.82	0.42
2:C:113:GLU:O	2:C:116:LEU:N	2.47	0.42
2:C:22:SER:N	2:C:25:GLU:OE2	2.53	0.41
2:D:49:TRP:CD2	2:D:57:PRO:HD2	2.56	0.41
2:C:187:SER:OG	2:C:190:HIS:HD2	2.02	0.41
1:A:3:GLU:O	1:A:5:ASP:N	2.53	0.41
2:C:19:THR:CG2	2:C:20:GLU:HG3	2.27	0.41
1:B:30:ILE:CG2	2:C:161:VAL:HG11	2.50	0.41
1:B:98:GLN:NE2	2:C:192:ARG:CG	2.84	0.41
1:A:21:LYS:HB3	1:A:21:LYS:HZ2	1.84	0.41
2:D:102:VAL:HG23	2:D:110:TYR:HB2	2.02	0.41
2:C:78:ARG:NH1	2:C:88:VAL:HG21	2.36	0.41
2:D:139:CYS:O	2:D:142:VAL:HG21	2.20	0.41
1:B:98:GLN:NE2	2:C:192:ARG:NE	2.49	0.41
2:D:10:ASN:ND2	2:D:12:ARG:HB2	2.36	0.41
2:C:7:GLN:HA	2:C:12:ARG:O	2.21	0.41
1:A:44:CYS:CB	1:A:79:SER:O	2.69	0.41
2:C:74:GLY:HA2	2:C:90:ALA:O	2.21	0.41
2:D:43:ASP:HB2	2:D:44:ASN:H	1.74	0.41
2:C:120:ARG:HG3	2:C:124:GLN:O	2.21	0.41
2:D:287:CYS:O	2:D:288:ASN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:104:GLY:C	2:D:106:ASP:N	2.74	0.41
1:B:46:SER:N	6:B:529:HOH:O	2.46	0.41
2:D:47:PHE:O	2:D:50:MET:HB2	2.21	0.41
1:A:95:ASP:C	1:A:95:ASP:OD1	2.59	0.41
1:B:77:LEU:HA	1:B:77:LEU:HD23	1.67	0.41
2:D:22:SER:HB3	2:D:25:GLU:HG3	2.01	0.41
2:D:213:ALA:O	2:D:229:TRP:HH2	2.04	0.41
2:D:200:ARG:CG	2:D:201:SER:N	2.82	0.40
1:A:87:ARG:HB2	1:A:111:GLU:OE1	2.21	0.40
1:B:101:ILE:HG22	1:B:102:LYS:N	2.35	0.40
1:B:4:CYS:SG	1:B:14:LYS:HG2	2.61	0.40
1:A:87:ARG:HG2	1:A:88:PRO:CD	2.49	0.40
2:D:246:PRO:O	2:D:247:ASP:C	2.60	0.40
1:A:89:MET:HG2	1:A:90:SER:N	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:233:VAL:O	2:C:235:ARG:NH1[2_765]	2.08	0.12

## 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	114/116 (98%)	105 (92%)	6 (5%)	3 (3%)	7	22
1	B	114/116 (98%)	99 (87%)	13 (11%)	2 (2%)	11	34
2	C	283/288 (98%)	251 (89%)	28 (10%)	4 (1%)	14	42
2	D	270/288 (94%)	239 (88%)	26 (10%)	5 (2%)	10	32
All	All	781/808 (97%)	694 (89%)	73 (9%)	14 (2%)	11	34

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	94	SER
2	D	43	ASP
2	D	148	SER
1	A	11	CYS
1	B	6	GLY
1	A	9	ASN
1	B	7	LYS
2	C	96	ILE
2	C	140	ARG
2	C	248	SER
2	D	122	LYS
2	D	15	VAL
2	C	15	VAL
1	A	44	CYS

### 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	97/102 (95%)	90 (93%)	7 (7%)	18	45
1	B	96/102 (94%)	93 (97%)	3 (3%)	47	81
2	C	225/248 (91%)	213 (95%)	12 (5%)	28	61
2	D	222/248 (90%)	212 (96%)	10 (4%)	34	68
All	All	640/700 (91%)	608 (95%)	32 (5%)	30	64

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	21	LYS
1	A	43	GLU
1	A	71	HIS
1	A	79	SER
1	A	90	SER

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Mol	Chain	Res	Type
1	A	102	LYS
1	B	77	LEU
1	B	102	LYS
1	B	106	GLN
2	C	15	VAL
2	C	37	THR
2	C	78	ARG
2	C	94	SER
2	C	114	CYS
2	C	116	LEU
2	C	191	LEU
2	C	222	THR
2	C	247	ASP
2	C	249	LYS
2	C	250	SER
2	C	254	VAL
2	D	14	GLN
2	D	18	LYS
2	D	37	THR
2	D	84	LYS
2	D	94	SER
2	D	120	ARG
2	D	141	ASP
2	D	156	ASN
2	D	165	ARG
2	D	191	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	26	ASN
1	A	98	GLN
1	B	98	GLN
1	B	99	ASN
2	C	14	GLN
2	C	80	ASN
2	C	130	GLN
2	C	132	GLN
2	C	190	HIS
2	C	259	ASN
2	D	7	GLN

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Mol	Chain	Res	Type
2	D	10	ASN
2	D	14	GLN
2	D	53	ASN
2	D	130	GLN
2	D	132	GLN
2	D	180	ASN
2	D	190	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, 5 are monoatomic - leaving 5 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	MLI	A	401	-	0,6,6	0.00	-	0,7,7	0.00	-
4	MLI	A	402	-	0,6,6	0.00	-	0,7,7	0.00	-
4	MLI	A	403	-	0,6,6	0.00	-	0,7,7	0.00	-
5	MPD	B	404	-	6,7,7	0.40	0	7,10,10	0.90	0
4	MLI	C	400	-	0,6,6	0.00	-	0,7,7	0.00	-

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	MLI	A	401	-	-	0/0/4/4	0/0/0/0
4	MLI	A	402	-	-	0/0/4/4	0/0/0/0
4	MLI	A	403	-	-	0/0/4/4	0/0/0/0
5	MPD	B	404	-	-	0/5/5/5	0/0/0/0
4	MLI	C	400	-	-	0/0/4/4	0/0/0/0

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.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	404	MPD	9	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, 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	116/116 (100%)	0.18	3 (2%) 59 47	43, 66, 93, 98	0
1	B	116/116 (100%)	0.21	3 (2%) 59 47	26, 56, 91, 99	0
2	C	287/288 (99%)	0.42	26 (9%) 11 6	23, 69, 100, 145	0
2	D	280/288 (97%)	0.57	33 (11%) 6 3	47, 82, 100, 102	0
All	All	799/808 (98%)	0.41	65 (8%) 15 7	23, 75, 99, 145	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	97	THR	7.4
2	D	85	PRO	7.2
2	C	85	PRO	7.1
2	C	171	ALA	5.5
2	C	17	TYR	5.1
2	C	69	VAL	5.1
2	D	170	PRO	5.0
2	C	86	ARG	4.6
2	C	21	LEU	4.6
1	B	2	LEU	4.5
2	D	263	ALA	4.3
2	D	251	ASP	4.3
2	C	172	SER	4.1
2	D	250	SER	4.0
2	D	83	ASN	3.9
2	D	228	LEU	3.8
2	D	246	PRO	3.7
2	D	172	SER	3.7
2	D	253	PRO	3.6
2	D	173	SER	3.5
2	D	93	CYS	3.5

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Mol	Chain	Res	Type	RSRZ
2	D	94	SER	3.5
2	D	254	VAL	3.5
2	D	110	TYR	3.4
2	D	131	TYR	3.4
2	D	89	CYS	3.3
2	C	28	SER	3.3
2	C	76	LYS	3.3
2	C	246	PRO	3.0
2	C	252	GLU	3.0
2	D	262	TYR	3.0
2	D	219	ILE	3.0
2	C	247	ASP	2.9
2	D	3	CYS	2.8
2	C	30	GLY	2.8
2	D	100	GLY	2.8
2	C	249	LYS	2.8
2	C	82	LYS	2.7
2	C	96	ILE	2.7
2	D	264	SER	2.7
2	C	22	SER	2.7
2	D	169	GLU	2.7
2	C	36	TRP	2.7
2	D	245	CYS	2.5
2	D	92	ASP	2.5
2	D	186	SER	2.5
1	A	116	SER	2.4
2	C	37	THR	2.4
1	A	2	LEU	2.4
2	D	91	PRO	2.4
2	C	78	ARG	2.3
1	B	116	SER	2.3
1	B	10	ILE	2.2
2	C	68	ASN	2.2
2	D	218	ASP	2.2
2	C	250	SER	2.2
2	D	174	GLU	2.2
2	D	115	ALA	2.2
2	C	93	CYS	2.2
2	D	217	GLU	2.1
2	D	84	LYS	2.1
2	C	81	LYS	2.1
1	A	1	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
2	D	73	PRO	2.0
2	C	1	GLY	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
4	MLI	C	400	7/7	0.53	0.55	7.30	75,75,76,76	7
5	MPD	B	404	8/8	0.93	0.51	6.40	71,71,72,73	0
4	MLI	A	403	7/7	0.88	0.40	6.13	44,46,49,49	7
4	MLI	A	402	7/7	0.80	0.40	4.94	53,53,55,56	7
3	IR3	C	505	1/1	0.77	1.11	-	97,97,97,97	1
3	IR3	B	502	1/1	0.98	0.32	-	79,79,79,79	1
4	MLI	A	401	7/7	0.79	0.31	-	64,65,66,66	7
3	IR3	C	503	1/1	0.91	0.27	-	94,94,94,94	1
3	IR3	A	504	1/1	0.84	0.97	-	70,70,70,70	1
3	IR3	D	501	1/1	0.86	0.52	-	94,94,94,94	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.