



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:04 PM GMT

PDB ID : 1WDL  
Title : fatty acid beta-oxidation multienzyme complex from *Pseudomonas fragi*, form II (native4)  
Authors : Ishikawa, M.; Tsuchiya, D.; Oyama, T.; Tsunaka, Y.; Morikawa, K.  
Deposited on : 2004-05-17  
Resolution : 3.50 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

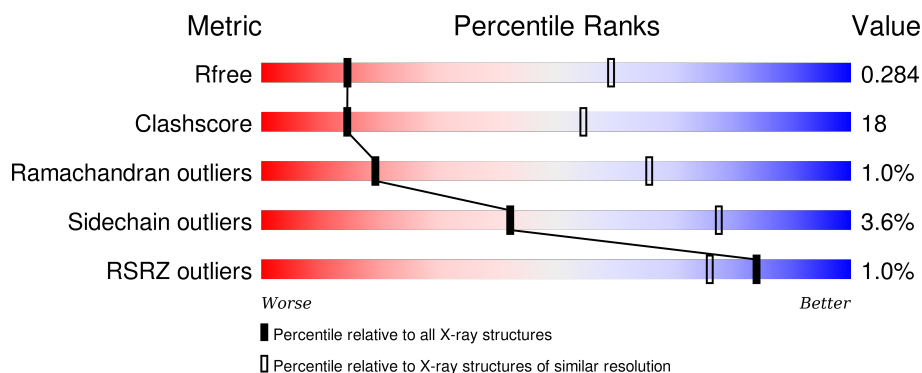
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	715	<div> <div> <div>0%</div> <div>63%</div> <div>34%</div> <div>•</div> </div> </div>
1	B	715	<div> <div>2%</div> <div>62%</div> <div>35%</div> <div>••</div> </div>
2	C	390	<div> <div>64%</div> <div>33%</div> <div>•</div> </div>
2	D	390	<div> <div>64%</div> <div>35%</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
3	ACO	C	3001	-	-	-	X
3	ACO	D	4001	-	-	-	X
5	N8E	A	1002	-	-	-	X
5	N8E	A	1004	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 16610 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.

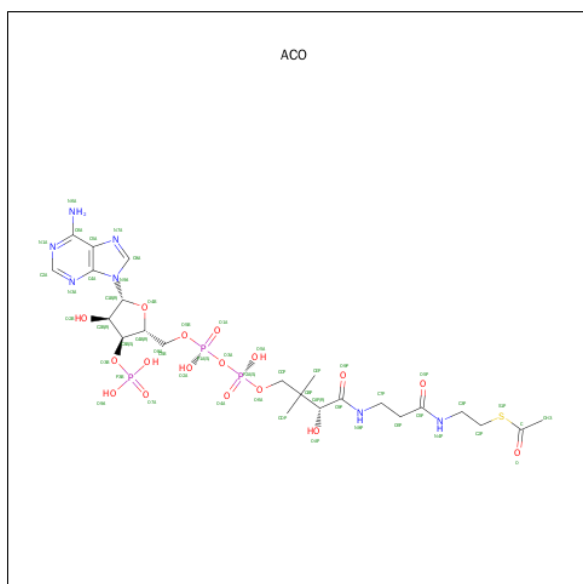
- Molecule 1 is a protein called Fatty oxidation complex alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	715	Total	C	N	O	S	0	0	0
			5361	3420	906	1008	27			
1	B	706	Total	C	N	O	S	0	0	0
			5241	3340	879	995	27			

- Molecule 2 is a protein called 3-ketoacyl-CoA thiolase.

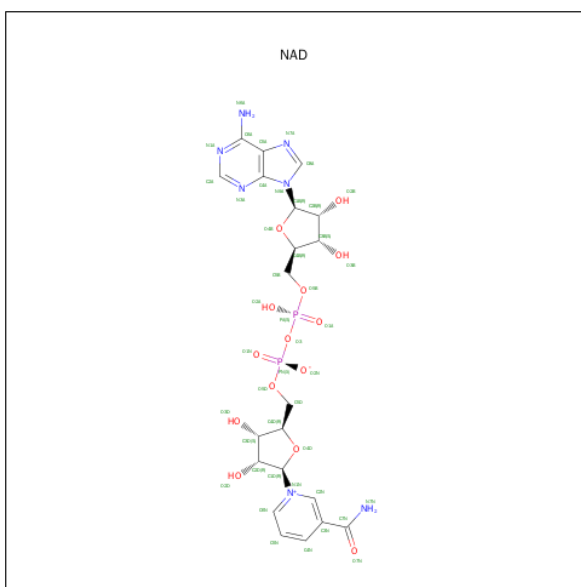
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	390	Total	C	N	O	S	0	0	0
			2885	1795	513	548	29			
2	D	390	Total	C	N	O	S	0	0	0
			2881	1792	512	548	29			

- Molecule 3 is ACETYL COENZYME \*A (three-letter code: ACO) (formula: C<sub>23</sub>H<sub>38</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>S).



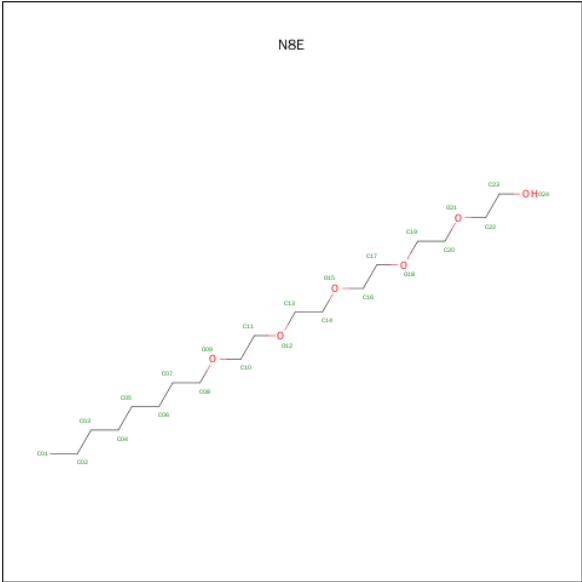
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	P	0	0
			40	16	6	15	3		
3	D	1	Total	C	N	O	P	S	0
			51	23	7	17	3	1	0

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			35	15	5	13	2		
4	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 5 is 3,6,9,12,15-PENTAOXATRICOSAN-1-OL (three-letter code: N8E) (formula:  $C_{18}H_{38}O_6$ ).

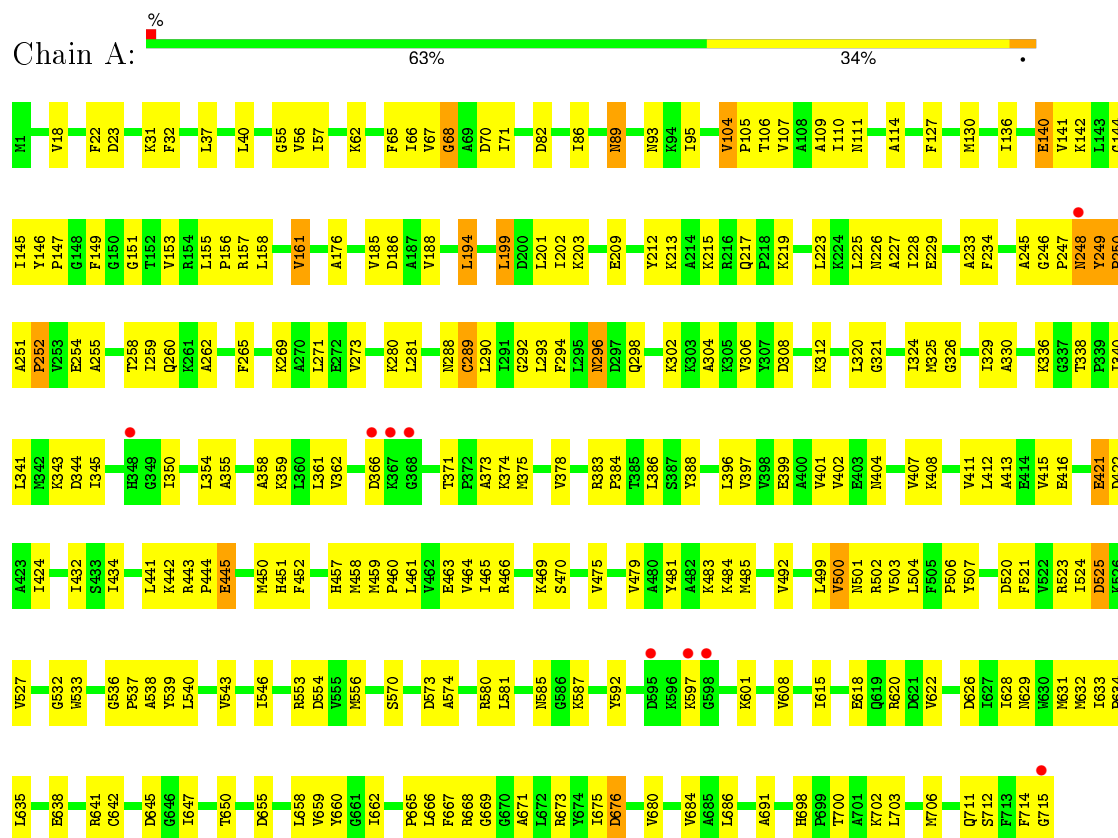


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			24	18	6		
5	A	1	Total	C	O	0	0
			24	18	6		
5	A	1	Total	C	O	0	0
			24	18	6		

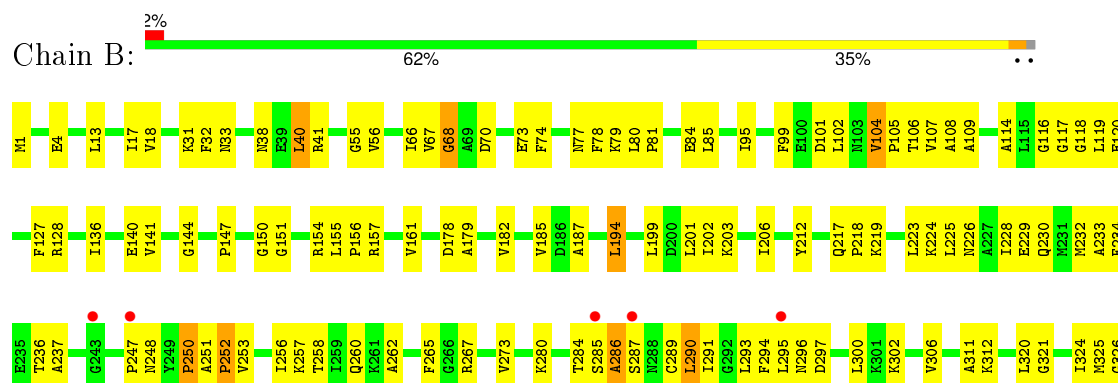
### 3 Residue-property plots

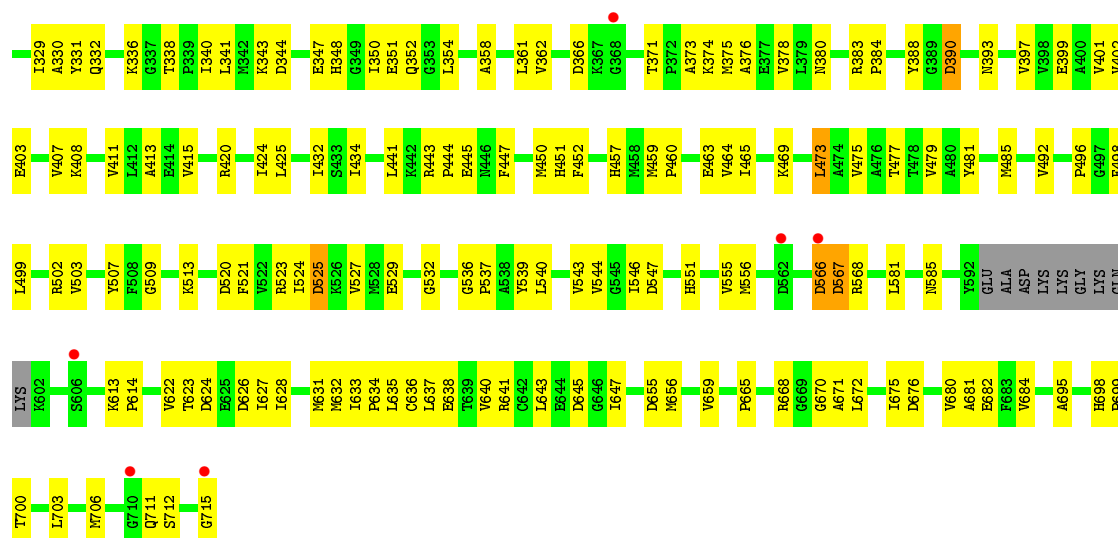
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Fatty oxidation complex alpha subunit



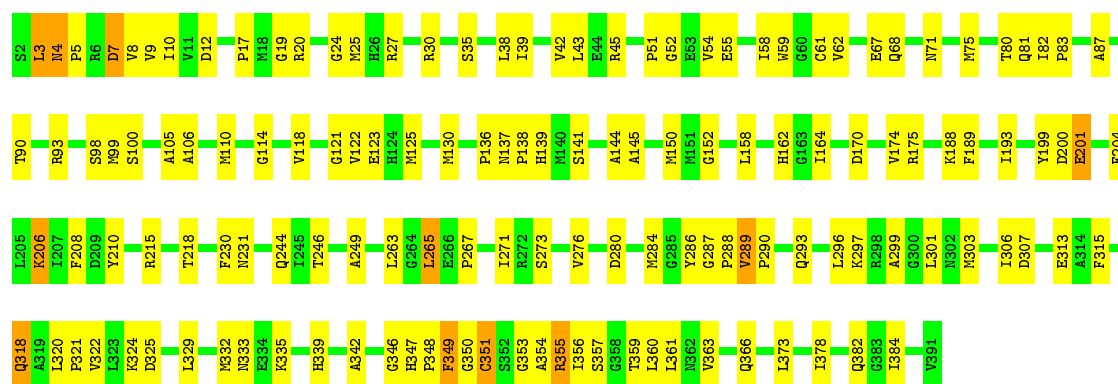
#### • Molecule 1: Fatty oxidation complex alpha subunit





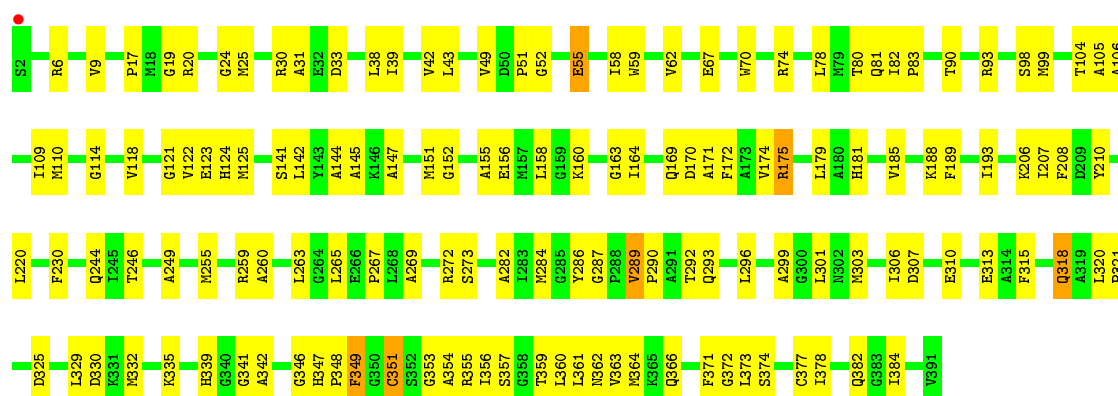
• Molecule 2: 3-ketoacyl-CoA thiolase

Chain C: 64% 33%



• Molecule 2: 3-ketoacyl-CoA thiolase

Chain D: 64% 35%





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.66Å 118.41Å 191.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 3.50 19.76 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (12.00-3.50) 99.6 (19.76-3.50)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.18 (at 3.52Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.233 , 0.283 0.237 , 0.284	Depositor DCC
$R_{free}$ test set	1875 reflections (6.88%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.9	Xtriage
Anisotropy	0.507	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 27.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 27822 reflections	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	16610	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: N8E, ACO, NAD

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	0.26	0/5448	0.53	1/7364 (0.0%)
1	B	0.25	0/5325	0.53	3/7213 (0.0%)
2	C	0.24	0/2933	0.50	0/3959
2	D	0.24	0/2929	0.49	0/3955
All	All	0.25	0/16635	0.52	4/22491 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	568	ARG	N-CA-C	-8.61	87.75	111.00
1	A	715	GLY	N-CA-C	-7.12	95.30	113.10
1	B	715	GLY	N-CA-C	-6.87	95.92	113.10
1	B	286	ALA	N-CA-C	-5.27	96.78	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	5361	0	5415	212	0
1	B	5241	0	5227	179	0
2	C	2885	0	2881	110	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2881	0	2870	115	0
3	C	40	0	21	0	0
3	D	51	0	34	4	0
4	A	35	0	19	0	0
4	B	44	0	26	1	0
5	A	72	0	114	6	0
All	All	16610	0	16607	596	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 18.

All (596) 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:289:VAL:HG13	2:C:290:PRO:HD3	1.49	0.95
2:D:272:ARG:HH12	2:D:301:LEU:HD21	1.33	0.94
2:D:289:VAL:HG13	2:D:290:PRO:HD3	1.49	0.93
1:A:666:LEU:H	1:A:666:LEU:HD12	1.35	0.90
2:D:378:ILE:HB	2:D:382:GLN:HB2	1.57	0.86
1:A:354:LEU:HD13	1:B:380:ASN:HB3	1.56	0.86
2:D:272:ARG:NH1	2:D:301:LEU:HD21	1.93	0.84
1:A:107:VAL:HG22	1:A:127:PHE:HB2	1.59	0.83
1:B:114:ALA:HB3	1:B:136:ILE:HG22	1.59	0.83
1:B:107:VAL:HG22	1:B:127:PHE:HB2	1.59	0.83
1:A:114:ALA:HB3	1:A:136:ILE:HG22	1.61	0.82
2:D:348:PRO:HB2	2:D:351:CYS:HB3	1.59	0.82
2:C:55:GLU:HG3	2:C:114:GLY:HA2	1.64	0.80
1:B:290:LEU:HG	1:B:293:LEU:HD12	1.63	0.80
1:B:320:LEU:HD11	1:B:415:VAL:HG21	1.66	0.78
1:B:144:GLY:HA2	1:B:250:PRO:HB3	1.65	0.77
2:D:263:LEU:HB2	2:D:265:LEU:HD22	1.66	0.77
2:C:306:ILE:HD13	2:C:373:LEU:HB2	1.67	0.77
1:A:465:ILE:H	1:A:465:ILE:HD12	1.49	0.76
1:B:247:PRO:HB3	1:B:529:GLU:O	1.84	0.76
1:B:56:VAL:HB	1:B:106:THR:HG22	1.68	0.76
1:B:99:PHE:HA	1:B:102:LEU:HD13	1.69	0.75
1:A:254:GLU:HG3	1:A:281:LEU:HD21	1.67	0.74
1:A:354:LEU:HD21	1:A:384:PRO:HG3	1.70	0.74
1:B:258:THR:HG23	1:B:273:VAL:HG12	1.71	0.73
1:B:361:LEU:HB3	1:B:375:MET:HG3	1.72	0.71
1:A:141:VAL:HG12	1:A:251:ALA:HB3	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:ASP:O	1:B:567:ASP:HB2	1.90	0.71
1:B:248:ASN:HB3	1:B:532:GLY:O	1.90	0.71
2:D:55:GLU:HG3	2:D:114:GLY:HA2	1.72	0.71
1:B:248:ASN:O	1:B:250:PRO:HD3	1.91	0.70
1:A:404:ASN:HD21	5:A:1004:N8E:H031	1.54	0.70
1:A:296:ASN:HD21	1:A:650:THR:HB	1.56	0.70
1:B:141:VAL:HG12	1:B:251:ALA:HB3	1.73	0.70
2:D:125:MET:HE2	2:D:348:PRO:HA	1.73	0.70
2:D:306:ILE:HD13	2:D:373:LEU:HB2	1.71	0.70
2:C:25:MET:HE1	2:C:208:PHE:HB2	1.74	0.70
1:B:680:VAL:HG21	1:B:712:SER:HA	1.73	0.69
1:A:336:LYS:HG3	1:A:485:MET:HA	1.74	0.69
2:C:123:GLU:HG2	2:C:349:PHE:HB2	1.73	0.69
1:A:361:LEU:HB3	1:A:375:MET:HG3	1.74	0.69
1:B:336:LYS:HG3	1:B:485:MET:HA	1.75	0.69
1:A:680:VAL:HG21	1:A:712:SER:HA	1.74	0.68
1:A:626:ASP:HA	1:A:629:ASN:HD22	1.59	0.68
2:C:348:PRO:HB2	2:C:351:CYS:HB3	1.75	0.68
1:B:465:ILE:HD12	1:B:465:ILE:H	1.59	0.68
1:A:581:LEU:HB2	1:A:585:ASN:HB2	1.77	0.67
2:C:137:ASN:HD21	2:C:139:HIS:HB2	1.59	0.67
2:C:263:LEU:HB2	2:C:265:LEU:HD22	1.76	0.67
2:C:62:VAL:HG21	2:C:349:PHE:HB3	1.77	0.67
1:B:280:LYS:HB2	1:B:280:LYS:NZ	2.10	0.67
1:A:504:LEU:HD12	1:A:662:ILE:HG22	1.76	0.66
2:D:62:VAL:HG21	2:D:349:PHE:HB3	1.78	0.66
1:B:262:ALA:HA	1:B:265:PHE:HD1	1.62	0.64
2:C:10:ILE:HD11	2:C:271:ILE:HG13	1.77	0.64
1:B:31:LYS:HD2	1:B:70:ASP:HB2	1.79	0.64
2:D:98:SER:HB3	2:D:354:ALA:H	1.63	0.64
1:B:633:ILE:HB	1:B:634:PRO:HD3	1.78	0.64
1:B:547:ASP:HB3	1:B:581:LEU:HD22	1.79	0.64
1:B:32:PHE:HE2	1:B:40:LEU:HD12	1.62	0.64
2:C:43:LEU:HD11	2:C:82:ILE:HD11	1.80	0.64
1:B:613:LYS:HB3	1:B:614:PRO:HD3	1.81	0.63
1:A:633:ILE:HB	1:A:634:PRO:HD3	1.79	0.63
1:B:706:MET:HB2	1:B:711:GLN:HB2	1.79	0.63
2:C:51:PRO:HB3	2:C:82:ILE:HD13	1.80	0.63
2:D:289:VAL:HG23	2:D:293:GLN:NE2	2.13	0.63
2:D:25:MET:HE3	2:D:208:PHE:HB2	1.80	0.63
1:A:89:ASN:N	1:A:89:ASN:HD22	1.95	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ILE:HD13	1:A:290:LEU:HD11	1.80	0.62
1:A:330:ALA:HA	1:A:340:ILE:HD13	1.80	0.62
2:C:289:VAL:O	2:C:293:GLN:HG3	1.99	0.62
1:A:500:VAL:HG13	1:A:662:ILE:HG23	1.81	0.62
2:C:320:LEU:HB2	2:C:321:PRO:HD3	1.80	0.62
1:A:338:THR:HG21	1:A:481:TYR:HE1	1.64	0.62
2:D:320:LEU:HB2	2:D:321:PRO:HD3	1.81	0.62
2:D:342:ALA:HB1	2:D:347:HIS:HB2	1.82	0.62
2:C:315:PHE:O	2:C:318:GLN:HG3	2.00	0.62
1:A:186:ASP:CG	1:A:215:LYS:NZ	2.53	0.62
2:C:93:ARG:NH1	2:C:276:VAL:HG11	2.15	0.62
2:C:201:GLU:H	2:C:201:GLU:CD	2.04	0.61
2:C:125:MET:HE2	2:C:348:PRO:HA	1.82	0.60
1:A:592:TYR:HA	1:A:601:LYS:O	2.01	0.60
1:B:354:LEU:HD21	1:B:384:PRO:HG3	1.84	0.60
1:B:388:TYR:CE2	1:B:415:VAL:HG22	2.37	0.59
2:C:3:LEU:HG	2:C:3:LEU:O	2.02	0.59
2:D:30:ARG:HB2	2:D:33:ASP:OD2	2.01	0.59
1:A:289:CYS:HB2	1:A:714:PHE:O	2.02	0.59
1:B:330:ALA:HA	1:B:340:ILE:HD13	1.83	0.59
2:D:188:LYS:HD2	2:D:366:GLN:OE1	2.03	0.59
2:D:272:ARG:HH11	2:D:371:PHE:HE2	1.50	0.59
2:C:296:LEU:HD21	2:C:306:ILE:HD11	1.85	0.59
1:B:393:ASN:HA	1:B:420:ARG:HH12	1.67	0.59
1:A:343:LYS:HD3	1:A:344:ASP:N	2.17	0.59
1:B:581:LEU:HB2	1:B:585:ASN:HB2	1.85	0.59
1:A:302:LYS:O	1:A:306:VAL:HG23	2.03	0.59
1:A:141:VAL:HG23	1:A:142:LYS:HD2	1.84	0.59
1:A:362:VAL:HG21	1:B:373:ALA:HA	1.85	0.59
1:B:521:PHE:HA	1:B:524:ILE:HD12	1.85	0.59
1:A:109:ALA:HB1	1:A:194:LEU:HD22	1.83	0.59
1:A:226:ASN:OD1	1:A:228:ILE:HG22	2.03	0.58
1:A:669:GLY:HA3	1:A:673:ARG:HD3	1.84	0.58
1:A:212:TYR:OH	1:A:213:LYS:HE2	2.04	0.58
1:B:311:ALA:HB1	1:B:477:THR:HA	1.84	0.58
1:A:533:TRP:HH2	1:A:662:ILE:HD12	1.69	0.58
2:D:313:GLU:CD	2:D:341:GLY:HA3	2.23	0.58
1:B:622:VAL:HG23	1:B:626:ASP:HB2	1.86	0.58
1:B:228:ILE:HG22	2:D:282:ALA:HB3	1.85	0.58
2:C:287:GLY:O	2:C:290:PRO:HD2	2.04	0.58
2:D:359:THR:O	2:D:363:VAL:HG23	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:378:ILE:HB	2:C:382:GLN:HB2	1.85	0.57
1:A:521:PHE:HA	1:A:524:ILE:HD12	1.86	0.57
1:A:233:ALA:HB2	2:C:145:ALA:HB2	1.85	0.57
2:D:287:GLY:O	2:D:290:PRO:HD2	2.04	0.57
1:B:320:LEU:HD13	1:B:411:VAL:HG12	1.87	0.57
1:A:525:ASP:HB3	1:A:536:GLY:HA3	1.86	0.57
1:B:33:ASN:HB2	1:B:73:GLU:OE1	2.05	0.57
1:A:144:GLY:HA2	1:A:250:PRO:HB3	1.86	0.57
1:B:388:TYR:HE2	1:B:415:VAL:HG22	1.68	0.57
2:D:39:ILE:HG23	2:D:118:VAL:HG11	1.87	0.57
2:D:377:CYS:SG	3:D:4001:ACO:HH31	2.44	0.57
1:A:361:LEU:HD21	1:A:378:VAL:HG11	1.86	0.56
1:B:343:LYS:HD3	1:B:344:ASP:N	2.20	0.56
1:B:638:GLU:HA	1:B:641:ARG:HH11	1.70	0.56
1:B:671:ALA:O	1:B:675:ILE:HG13	2.05	0.56
1:A:633:ILE:HD11	1:A:686:LEU:HD22	1.88	0.56
1:B:424:ILE:HD12	1:B:424:ILE:N	2.21	0.56
1:B:109:ALA:HB1	1:B:194:LEU:HD22	1.85	0.56
1:B:217:GLN:HB3	1:B:218:PRO:HD3	1.87	0.56
1:A:226:ASN:CG	1:A:228:ILE:HG22	2.26	0.56
2:C:342:ALA:HB1	2:C:347:HIS:HB2	1.86	0.56
1:A:290:LEU:HD23	1:A:290:LEU:O	2.06	0.56
2:C:52:GLY:HA2	2:C:83:PRO:HG2	1.88	0.56
1:B:434:ILE:HD13	1:B:465:ILE:HG21	1.87	0.56
2:D:356:ILE:HG13	2:D:357:SER:N	2.21	0.55
1:B:67:VAL:O	1:B:68:GLY:O	2.23	0.55
2:D:315:PHE:O	2:D:318:GLN:HG3	2.06	0.55
2:C:130:MET:O	2:D:70:TRP:HZ2	1.89	0.55
1:A:31:LYS:HD2	1:A:70:ASP:HB2	1.87	0.55
1:A:209:GLU:OE2	2:D:206:LYS:HB2	2.07	0.55
1:A:466:ARG:HG3	1:A:470:SER:HB2	1.87	0.55
2:D:51:PRO:HB3	2:D:82:ILE:HD13	1.86	0.55
1:B:525:ASP:HB2	1:B:536:GLY:HA3	1.88	0.55
1:A:465:ILE:N	1:A:465:ILE:HD12	2.19	0.55
1:B:347:GLU:O	1:B:351:GLU:HG2	2.06	0.55
2:C:152:GLY:HA3	2:C:230:PHE:CD2	2.41	0.55
2:C:80:THR:HG22	2:C:81:GLN:N	2.22	0.55
1:A:451:HIS:HB3	1:A:463:GLU:HB2	1.89	0.55
1:B:150:GLY:HA3	1:B:154:ARG:NH1	2.23	0.54
1:A:450:MET:HG2	1:A:452:PHE:CE1	2.42	0.54
1:B:361:LEU:HD21	1:B:378:VAL:HG11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:397:VAL:HB	1:B:425:LEU:HD23	1.90	0.54
1:A:228:ILE:HG21	2:C:280:ASP:CG	2.27	0.54
2:C:303:MET:HE3	2:C:306:ILE:HD12	1.88	0.54
1:A:341:LEU:HD23	1:A:383:ARG:HB2	1.89	0.54
2:D:172:PHE:HE2	2:D:313:GLU:HG2	1.72	0.54
1:B:155:LEU:HB3	1:B:156:PRO:HD3	1.90	0.54
1:A:246:GLY:HA3	1:A:668:ARG:HH22	1.73	0.54
1:A:502:ARG:HD3	1:A:642:CYS:SG	2.48	0.54
1:B:413:ALA:HA	1:B:441:LEU:HD23	1.91	0.53
1:A:503:VAL:HG12	1:A:635:LEU:HG	1.89	0.53
2:D:156:GLU:O	2:D:160:LYS:HG2	2.08	0.53
1:A:186:ASP:CG	1:A:215:LYS:HZ3	2.12	0.53
1:A:153:VAL:HG21	1:A:271:LEU:HD23	1.91	0.53
1:A:539:TYR:O	1:A:543:VAL:HG23	2.08	0.53
2:C:359:THR:O	2:C:363:VAL:HG23	2.07	0.53
1:A:434:ILE:HD13	1:A:465:ILE:HG21	1.89	0.53
1:A:706:MET:HB2	1:A:711:GLN:HB2	1.90	0.53
2:C:4:ASN:HB2	2:C:7:ASP:OD2	2.09	0.53
1:B:229:GLU:O	2:D:145:ALA:HB2	2.09	0.53
2:D:272:ARG:NH1	2:D:371:PHE:HE2	2.08	0.52
1:A:292:GLY:O	1:A:296:ASN:HB2	2.09	0.52
2:D:80:THR:HG22	2:D:81:GLN:N	2.23	0.52
1:A:457:HIS:CE1	1:A:458:MET:HG2	2.44	0.52
1:A:258:THR:HG23	1:A:273:VAL:HG12	1.90	0.52
1:B:117:GLY:HA2	1:B:120:GLU:OE1	2.09	0.52
2:D:151:MET:HE3	3:D:4001:ACO:HH32	1.91	0.52
1:B:450:MET:HG2	1:B:452:PHE:CE1	2.45	0.52
2:C:164:ILE:HD12	2:C:164:ILE:N	2.24	0.52
1:A:248:ASN:O	1:A:250:PRO:HD3	2.08	0.52
1:B:507:TYR:HB2	1:B:635:LEU:HD12	1.91	0.52
2:D:175:ARG:HE	2:D:179:LEU:HD12	1.74	0.52
2:D:106:ALA:O	2:D:110:MET:HG3	2.10	0.52
1:B:202:ILE:O	1:B:206:ILE:HG13	2.10	0.52
1:A:32:PHE:CE1	1:A:66:ILE:HG21	2.44	0.52
2:D:246:THR:OG1	2:D:346:GLY:HA3	2.10	0.52
1:A:251:ALA:N	1:A:252:PRO:CD	2.73	0.52
2:D:38:LEU:O	2:D:42:VAL:HG23	2.10	0.52
1:A:265:PHE:HB3	1:A:269:LYS:HB3	1.90	0.52
1:A:671:ALA:O	1:A:675:ILE:HG13	2.09	0.52
2:C:246:THR:OG1	2:C:346:GLY:HA3	2.10	0.52
1:B:325:MET:HG3	4:B:2001:NAD:O4D	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:348:HIS:O	1:B:352:GLN:HG2	2.10	0.52
2:D:189:PHE:HB3	2:D:193:ILE:CD1	2.40	0.52
1:A:155:LEU:HB3	1:A:156:PRO:HD3	1.92	0.52
1:B:523:ARG:O	1:B:527:VAL:HG23	2.09	0.52
1:A:525:ASP:CB	1:A:536:GLY:HA3	2.40	0.51
2:C:20:ARG:O	2:C:24:GLY:HA3	2.10	0.51
2:C:10:ILE:N	2:C:10:ILE:HD12	2.25	0.51
1:B:330:ALA:HA	1:B:340:ILE:HG21	1.93	0.51
1:B:393:ASN:HA	1:B:420:ARG:NH1	2.26	0.51
2:D:122:VAL:HG23	2:D:249:ALA:HB2	1.91	0.51
1:B:475:VAL:O	1:B:479:VAL:HG23	2.10	0.51
1:A:312:LYS:NZ	1:A:422:ASP:HB2	2.25	0.51
1:B:280:LYS:HB2	1:B:280:LYS:HZ3	1.76	0.51
2:C:318:GLN:C	2:C:321:PRO:HD2	2.31	0.51
1:A:292:GLY:HA3	1:A:714:PHE:O	2.11	0.51
2:C:99:MET:HB3	2:C:384:ILE:HD13	1.93	0.51
1:B:141:VAL:HG12	1:B:251:ALA:CB	2.39	0.50
2:D:80:THR:HG22	2:D:81:GLN:H	1.76	0.50
1:A:507:TYR:HB2	1:A:635:LEU:HD12	1.94	0.50
1:B:525:ASP:CB	1:B:536:GLY:HA3	2.42	0.50
2:D:9:VAL:HG13	2:D:267:PRO:HB3	1.94	0.50
2:D:310:GLU:HG3	2:D:360:LEU:HB2	1.93	0.50
1:A:523:ARG:O	1:A:527:VAL:HG23	2.12	0.50
1:B:262:ALA:HA	1:B:265:PHE:CD1	2.44	0.50
2:D:313:GLU:O	2:D:342:ALA:HB3	2.11	0.50
2:D:265:LEU:HD23	2:D:265:LEU:H	1.76	0.50
1:A:228:ILE:HD13	2:C:280:ASP:HB2	1.93	0.50
1:A:645:ASP:HB2	1:A:647:ILE:HD13	1.94	0.50
1:A:199:LEU:HD11	1:A:203:LYS:HE3	1.94	0.50
2:D:164:ILE:N	2:D:164:ILE:HD12	2.27	0.50
2:D:17:PRO:HB3	2:D:210:TYR:O	2.11	0.50
1:B:312:LYS:HG3	1:B:473:LEU:HD21	1.92	0.50
1:A:246:GLY:HA3	1:A:668:ARG:HH12	1.76	0.50
1:A:638:GLU:HA	1:A:641:ARG:HH11	1.76	0.50
1:A:157:ARG:HD2	1:A:219:LYS:HA	1.93	0.50
2:D:58:ILE:HD12	2:D:105:ALA:HB2	1.93	0.50
1:A:255:ALA:O	1:A:259:ILE:HG13	2.12	0.49
1:B:287:SER:OG	1:B:668:ARG:NH1	2.45	0.49
1:A:424:ILE:HD12	1:A:424:ILE:N	2.27	0.49
1:A:82:ASP:O	1:A:86:ILE:HG13	2.12	0.49
1:A:350:ILE:HG21	1:A:384:PRO:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:98:SER:CB	2:D:354:ALA:H	2.25	0.49
1:A:580:ARG:CZ	1:A:608:VAL:HG22	2.41	0.49
1:B:401:VAL:HG12	1:B:402:VAL:N	2.27	0.49
1:A:292:GLY:HA3	1:A:714:PHE:C	2.32	0.49
2:C:8:VAL:HG11	2:C:271:ILE:HD12	1.94	0.49
1:A:574:ALA:HB2	1:A:615:ILE:HD11	1.94	0.49
1:B:302:LYS:O	1:B:306:VAL:HG23	2.13	0.49
1:B:108:ALA:HB3	1:B:128:ARG:HG2	1.93	0.49
1:B:465:ILE:N	1:B:465:ILE:HD12	2.25	0.49
2:C:273:SER:HB3	2:C:299:ALA:HB2	1.94	0.49
1:B:199:LEU:HD11	1:B:203:LYS:HE3	1.92	0.49
1:B:101:ASP:CG	1:B:267:ARG:HH22	2.15	0.49
2:C:55:GLU:HG3	2:C:114:GLY:CA	2.40	0.49
1:A:325:MET:O	1:A:329:ILE:HG13	2.11	0.49
1:B:399:GLU:HG2	1:B:401:VAL:HG23	1.93	0.49
1:B:520:ASP:O	1:B:524:ILE:HG13	2.12	0.49
2:D:20:ARG:O	2:D:24:GLY:HA3	2.12	0.49
1:A:413:ALA:HA	1:A:441:LEU:HD23	1.93	0.49
1:A:371:THR:HG22	1:A:373:ALA:H	1.76	0.49
2:D:19:GLY:O	2:D:246:THR:HG23	2.13	0.49
1:A:465:ILE:H	1:A:465:ILE:CD1	2.24	0.49
1:B:465:ILE:HA	1:B:492:VAL:O	2.12	0.49
2:D:318:GLN:C	2:D:321:PRO:HD2	2.33	0.49
1:B:104:VAL:HG22	1:B:105:PRO:HD2	1.95	0.49
1:B:655:ASP:O	1:B:659:VAL:HG23	2.11	0.49
1:B:161:VAL:HG21	1:B:225:LEU:HD22	1.94	0.49
1:A:288:ASN:O	1:A:289:CYS:SG	2.68	0.49
2:C:355:ARG:NH2	2:C:356:ILE:HG22	2.28	0.49
1:B:289:CYS:C	1:B:291:ILE:H	2.11	0.49
1:A:404:ASN:HD21	5:A:1004:N8E:C03	2.25	0.49
1:B:622:VAL:CG2	1:B:626:ASP:HB2	2.42	0.49
2:D:170:ASP:O	2:D:174:VAL:HG23	2.13	0.49
1:A:553:ARG:HH12	1:A:573:ASP:CG	2.15	0.49
2:D:125:MET:CE	2:D:246:THR:H	2.26	0.48
2:C:17:PRO:HB2	2:C:25:MET:HE2	1.95	0.48
2:C:17:PRO:HB3	2:C:210:TYR:O	2.13	0.48
2:D:377:CYS:SG	3:D:4001:ACO:CH3	3.01	0.48
1:B:503:VAL:HG12	1:B:635:LEU:HG	1.95	0.48
1:A:537:PRO:HG2	1:A:631:MET:CE	2.43	0.48
1:A:443:ARG:N	1:A:444:PRO:HD3	2.28	0.48
2:D:244:GLN:HB2	2:D:346:GLY:HA2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:700:THR:OG1	1:A:703:LEU:HB2	2.13	0.48
1:A:658:LEU:HD22	1:A:662:ILE:HD13	1.94	0.48
2:C:9:VAL:HG13	2:C:267:PRO:HB3	1.94	0.48
2:C:158:LEU:O	2:C:158:LEU:HD23	2.13	0.48
1:A:228:ILE:HG23	1:A:229:GLU:N	2.27	0.48
2:C:313:GLU:O	2:C:342:ALA:HB3	2.13	0.48
1:A:499:LEU:O	1:A:503:VAL:HG23	2.13	0.48
2:C:162:HIS:HB3	2:C:324:LYS:HZ1	1.78	0.48
2:C:38:LEU:O	2:C:42:VAL:HG23	2.13	0.48
1:B:157:ARG:HG2	1:B:223:LEU:HD23	1.96	0.48
1:B:331:TYR:HE1	1:B:361:LEU:HG	1.79	0.48
1:A:358:ALA:O	1:A:362:VAL:HG23	2.14	0.48
2:C:289:VAL:CG1	2:C:290:PRO:HD3	2.33	0.48
1:A:111:ASN:HA	1:A:194:LEU:HD11	1.95	0.48
1:A:580:ARG:HG2	1:A:587:LYS:HB3	1.96	0.48
1:A:130:MET:SD	1:A:185:VAL:HG21	2.53	0.48
1:A:293:LEU:HD23	1:A:714:PHE:CE2	2.48	0.48
2:C:30:ARG:NH1	2:C:67:GLU:HG2	2.29	0.48
1:A:451:HIS:HB2	1:A:465:ILE:HD11	1.96	0.47
1:A:676:ASP:OD2	1:A:712:SER:HB3	2.14	0.47
1:B:338:THR:HG21	1:B:481:TYR:HE1	1.79	0.47
1:B:445:GLU:H	1:B:445:GLU:CD	2.16	0.47
2:D:357:SER:O	2:D:360:LEU:HB3	2.14	0.47
1:A:464:VAL:HG11	1:A:475:VAL:HG13	1.96	0.47
1:B:695:ALA:HA	1:B:698:HIS:ND1	2.29	0.47
2:C:8:VAL:HG21	2:C:106:ALA:HB1	1.97	0.47
2:D:289:VAL:CG1	2:D:290:PRO:HD3	2.33	0.47
1:B:294:PHE:HA	1:B:297:ASP:HB3	1.97	0.47
2:C:98:SER:HB3	2:C:353:GLY:HA3	1.96	0.47
2:D:307:ASP:O	2:D:335:LYS:HB2	2.14	0.47
1:B:41:ARG:HD2	1:B:95:ILE:HG12	1.95	0.47
2:C:200:ASP:OD2	2:C:204:PHE:HB2	2.15	0.47
2:D:164:ILE:HG22	2:D:169:GLN:HG3	1.97	0.47
1:A:110:ILE:HB	1:A:130:MET:HG3	1.95	0.47
1:A:502:ARG:NH1	1:A:647:ILE:HG12	2.30	0.47
2:C:9:VAL:CG1	2:C:267:PRO:HB3	2.45	0.47
2:D:360:LEU:HD22	2:D:374:SER:HB3	1.97	0.47
1:B:116:GLY:HA2	1:B:140:GLU:OE1	2.14	0.47
1:B:624:ASP:O	1:B:628:ILE:HG13	2.14	0.47
1:B:350:ILE:HG21	1:B:384:PRO:HB3	1.95	0.47
1:B:623:THR:O	1:B:627:ILE:HG13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:618:GLU:OE2	1:A:620:ARG:HD2	2.15	0.47
2:D:99:MET:HB3	2:D:384:ILE:HD13	1.95	0.47
1:A:289:CYS:SG	1:A:714:PHE:O	2.73	0.47
1:A:320:LEU:HD13	1:A:411:VAL:HG12	1.96	0.47
1:A:22:PHE:CD2	1:A:66:ILE:HD11	2.49	0.47
1:B:105:PRO:HG3	1:B:212:TYR:CD2	2.50	0.47
1:A:249:TYR:HD1	1:A:249:TYR:O	1.98	0.47
1:A:666:LEU:HD12	1:A:666:LEU:N	2.18	0.46
1:B:341:LEU:HD23	1:B:383:ARG:HB2	1.97	0.46
2:C:137:ASN:ND2	2:C:139:HIS:HB2	2.27	0.46
2:C:93:ARG:HG2	2:C:378:ILE:HD13	1.96	0.46
1:A:228:ILE:HD13	2:C:280:ASP:CB	2.45	0.46
1:A:506:PRO:HG2	1:A:638:GLU:HG2	1.97	0.46
1:A:421:GLU:HA	1:A:443:ARG:CZ	2.45	0.46
1:A:130:MET:HB2	1:A:185:VAL:HG11	1.97	0.46
1:A:416:GLU:OE1	1:A:442:LYS:HG2	2.14	0.46
1:A:23:ASP:OD2	1:A:62:LYS:HE2	2.15	0.46
1:B:253:VAL:O	1:B:257:LYS:HG3	2.15	0.46
1:A:304:ALA:O	1:A:308:ASP:HB2	2.15	0.46
1:A:401:VAL:HG12	1:A:402:VAL:N	2.30	0.46
2:C:208:PHE:HD2	2:C:210:TYR:HH	1.61	0.46
2:D:98:SER:HB2	2:D:353:GLY:HA3	1.97	0.46
1:A:140:GLU:HG3	1:A:147:PRO:HD3	1.97	0.46
2:D:59:TRP:HE1	2:D:122:VAL:HB	1.80	0.46
2:D:74:ARG:NH2	2:D:78:LEU:HD21	2.30	0.46
1:A:37:LEU:HD22	1:A:95:ILE:HD12	1.95	0.46
1:B:251:ALA:N	1:B:252:PRO:CD	2.78	0.46
1:B:150:GLY:HA3	1:B:154:ARG:HH12	1.81	0.46
2:D:175:ARG:NE	2:D:179:LEU:HD12	2.30	0.46
1:A:459:MET:HA	1:A:460:PRO:HD3	1.74	0.46
1:B:79:LYS:HA	1:B:284:THR:HG23	1.97	0.46
2:C:303:MET:CE	2:C:306:ILE:HD12	2.44	0.46
2:C:17:PRO:HB2	2:C:25:MET:CE	2.45	0.46
1:A:18:VAL:CG1	1:A:56:VAL:HG22	2.46	0.46
1:A:461:LEU:HB2	1:A:660:TYR:HB3	1.97	0.46
2:D:272:ARG:NH1	2:D:371:PHE:CE2	2.83	0.46
1:B:411:VAL:O	1:B:415:VAL:HG23	2.15	0.46
1:A:186:ASP:CG	1:A:215:LYS:HZ1	2.18	0.46
1:A:209:GLU:O	2:D:207:ILE:HD12	2.15	0.46
1:B:358:ALA:O	1:B:362:VAL:HG23	2.16	0.46
1:B:443:ARG:N	1:B:444:PRO:HD3	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:VAL:O	1:A:411:VAL:HG23	2.15	0.46
2:C:244:GLN:HB2	2:C:346:GLY:HA2	1.97	0.46
1:A:507:TYR:CE1	1:A:631:MET:HB3	2.51	0.46
2:C:200:ASP:HB2	2:C:201:GLU:OE2	2.14	0.46
2:C:93:ARG:HH12	2:C:276:VAL:HG11	1.80	0.46
1:A:246:GLY:HA3	1:A:668:ARG:NH2	2.31	0.46
2:C:4:ASN:HD22	2:C:4:ASN:HA	1.58	0.46
1:B:81:PRO:HD2	1:B:84:GLU:HB2	1.96	0.46
1:A:465:ILE:HA	1:A:492:VAL:O	2.16	0.45
2:C:80:THR:HG22	2:C:81:GLN:H	1.80	0.45
1:B:147:PRO:HB3	1:B:151:GLY:HA3	1.98	0.45
1:A:355:ALA:O	1:A:359:LYS:HG3	2.16	0.45
1:B:407:VAL:O	1:B:411:VAL:HG23	2.15	0.45
2:D:9:VAL:CG1	2:D:267:PRO:HB3	2.46	0.45
1:A:374:LYS:O	1:A:378:VAL:HG23	2.16	0.45
1:B:66:ILE:O	1:B:118:GLY:HA3	2.16	0.45
2:C:12:ASP:CG	2:C:45:ARG:HH21	2.19	0.45
1:A:691:ALA:HB1	1:A:698:HIS:CE1	2.52	0.45
1:B:320:LEU:HB2	1:B:399:GLU:HA	1.97	0.45
1:A:321:GLY:O	1:A:326:GLY:HA3	2.17	0.45
1:A:520:ASP:O	1:A:524:ILE:HG13	2.16	0.45
1:A:157:ARG:HH11	1:A:157:ARG:HG2	1.81	0.45
2:C:87:ALA:HB1	2:D:93:ARG:HH21	1.81	0.45
2:D:364:MET:SD	2:D:372:GLY:HA3	2.57	0.45
1:B:136:ILE:HD11	1:B:179:ALA:HB1	1.98	0.45
1:A:248:ASN:HD22	1:A:248:ASN:HA	1.49	0.45
1:B:665:PRO:HB2	1:B:668:ARG:HD2	1.98	0.45
1:A:396:LEU:HD23	1:A:397:VAL:N	2.31	0.45
1:A:501:ASN:HD22	5:A:1002:N8E:H013	1.82	0.45
1:B:502:ARG:HG3	1:B:556:MET:HE1	1.99	0.45
1:B:13:LEU:HB2	1:B:17:ILE:O	2.17	0.45
2:C:188:LYS:HD2	2:C:366:GLN:OE1	2.17	0.45
1:A:262:ALA:HB2	1:A:273:VAL:HG21	1.99	0.45
1:B:232:MET:O	1:B:236:THR:HG23	2.17	0.45
2:C:329:LEU:HD23	2:C:329:LEU:O	2.17	0.45
2:D:52:GLY:HA2	2:D:83:PRO:HG2	1.99	0.45
1:A:666:LEU:H	1:A:666:LEU:CD1	2.12	0.45
1:A:89:ASN:N	1:A:89:ASN:ND2	2.63	0.45
1:B:637:LEU:O	1:B:641:ARG:HG3	2.17	0.45
1:A:502:ARG:HH11	1:A:647:ILE:HG12	1.82	0.45
1:A:56:VAL:HB	1:A:106:THR:HG22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:71:ASN:O	2:C:75:MET:HG2	2.17	0.45
1:B:374:LYS:O	1:B:378:VAL:HG23	2.17	0.44
1:A:147:PRO:HB3	1:A:151:GLY:HA3	1.98	0.44
1:A:412:LEU:HB3	1:A:441:LEU:HD21	1.99	0.44
2:C:59:TRP:O	2:C:90:THR:HA	2.17	0.44
1:B:300:LEU:HD11	1:B:656:MET:HG3	1.98	0.44
2:D:171:ALA:HA	2:D:220:LEU:HD21	1.98	0.44
2:D:289:VAL:HG23	2:D:293:GLN:HE21	1.79	0.44
2:D:98:SER:OG	2:D:121:GLY:HA3	2.16	0.44
2:D:359:THR:HA	2:D:362:ASN:HD22	1.81	0.44
2:D:255:MET:HE3	2:D:259:ARG:HG3	1.99	0.44
1:B:459:MET:HA	1:B:460:PRO:HD3	1.75	0.44
1:A:401:VAL:HG21	1:A:411:VAL:HG21	1.98	0.44
1:B:233:ALA:HB2	2:D:145:ALA:HB2	1.98	0.44
2:D:125:MET:HE3	2:D:246:THR:H	1.82	0.44
2:C:106:ALA:O	2:C:110:MET:HG3	2.18	0.44
1:B:636:CYS:O	1:B:640:VAL:HG23	2.18	0.44
2:D:299:ALA:O	2:D:301:LEU:HG	2.18	0.44
1:A:537:PRO:HG2	1:A:631:MET:HE1	1.99	0.44
1:B:234:PHE:CZ	1:B:260:GLN:HA	2.53	0.44
2:C:175:ARG:NH1	2:C:333:ASN:OD1	2.50	0.44
2:C:286:TYR:O	2:C:289:VAL:HG13	2.17	0.44
1:A:622:VAL:HG23	1:A:626:ASP:HB2	1.98	0.44
2:C:125:MET:HE3	2:C:246:THR:H	1.82	0.44
2:C:265:LEU:HD23	2:C:265:LEU:H	1.83	0.44
1:A:570:SER:HA	1:A:615:ILE:HG23	2.00	0.44
2:C:273:SER:CB	2:C:299:ALA:HB2	2.47	0.44
1:B:237:ALA:HB1	1:B:256:ILE:HD13	1.99	0.44
1:B:332:GLN:OE1	1:B:457:HIS:HA	2.17	0.44
1:A:655:ASP:O	1:A:659:VAL:HG23	2.18	0.44
2:D:152:GLY:HA3	2:D:230:PHE:CD2	2.53	0.44
2:C:35:SER:O	2:C:39:ILE:HG13	2.17	0.44
1:A:445:GLU:HA	1:A:469:LYS:O	2.17	0.44
1:B:80:LEU:HB3	1:B:81:PRO:CD	2.47	0.44
1:A:161:VAL:HG21	1:A:225:LEU:HD22	2.00	0.44
1:A:684:VAL:HG22	1:A:703:LEU:HD22	2.00	0.44
1:A:523:ARG:HH22	1:A:622:VAL:HG13	1.83	0.44
2:C:288:PRO:HB2	2:C:322:VAL:HG11	1.99	0.44
2:C:318:GLN:O	2:C:322:VAL:HG23	2.18	0.44
1:A:234:PHE:CZ	1:A:260:GLN:HA	2.53	0.44
2:D:163:GLY:C	2:D:164:ILE:HD12	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:LEU:HD23	1:B:294:PHE:HE1	1.83	0.43
1:A:289:CYS:CB	1:A:714:PHE:O	2.64	0.43
1:A:680:VAL:O	1:A:684:VAL:HG23	2.19	0.43
1:B:499:LEU:O	1:B:503:VAL:HG23	2.18	0.43
2:D:59:TRP:O	2:D:90:THR:HA	2.18	0.43
2:C:98:SER:HB3	2:C:353:GLY:CA	2.48	0.43
2:D:255:MET:HE1	2:D:260:ALA:HA	1.99	0.43
1:B:232:MET:HG3	2:D:147:ALA:HB3	2.00	0.43
1:B:321:GLY:O	1:B:326:GLY:HA3	2.17	0.43
1:B:551:HIS:O	1:B:555:VAL:HG23	2.18	0.43
2:D:286:TYR:O	2:D:289:VAL:HG13	2.18	0.43
2:D:303:MET:HE3	2:D:306:ILE:HD12	1.99	0.43
1:A:246:GLY:HA3	1:A:668:ARG:NH1	2.33	0.43
1:B:624:ASP:HA	1:B:627:ILE:HD12	2.00	0.43
2:C:141:SER:HA	2:C:144:ALA:O	2.18	0.43
2:C:122:VAL:HG23	2:C:249:ALA:HB2	2.00	0.43
2:C:136:PRO:O	2:C:138:PRO:HD3	2.18	0.43
1:B:226:ASN:HB3	1:B:229:GLU:H	1.84	0.43
1:A:71:ILE:O	1:A:71:ILE:HG22	2.19	0.43
1:B:465:ILE:HG12	1:B:498:PHE:CE2	2.54	0.43
2:C:93:ARG:HD2	2:C:100:SER:OG	2.19	0.43
1:A:503:VAL:C	1:A:506:PRO:HD2	2.38	0.43
2:C:339:HIS:CD2	2:C:363:VAL:HG22	2.54	0.43
2:C:170:ASP:O	2:C:174:VAL:HG23	2.17	0.43
1:A:245:ALA:HA	1:A:667:PHE:HB2	2.01	0.43
1:B:684:VAL:HG13	1:B:699:PRO:HG2	2.00	0.43
1:A:350:ILE:CG2	1:A:384:PRO:HB3	2.49	0.43
1:B:465:ILE:HG12	1:B:498:PHE:CZ	2.54	0.43
2:D:123:GLU:HG2	2:D:349:PHE:HB2	2.00	0.43
1:A:502:ARG:HA	1:A:556:MET:CE	2.49	0.43
1:B:537:PRO:HG2	1:B:631:MET:CE	2.48	0.43
1:B:546:ILE:N	1:B:546:ILE:HD12	2.34	0.43
2:C:58:ILE:HD12	2:C:105:ALA:HB2	2.00	0.43
1:A:622:VAL:CG2	1:A:626:ASP:HB2	2.49	0.43
2:D:141:SER:HA	2:D:144:ALA:O	2.19	0.43
1:B:645:ASP:HB2	1:B:647:ILE:HD13	2.00	0.43
1:B:296:ASN:N	1:B:296:ASN:HD22	2.15	0.43
2:C:19:GLY:O	2:C:246:THR:HG23	2.18	0.43
2:D:98:SER:HB2	2:D:353:GLY:CA	2.49	0.43
2:D:98:SER:HB3	2:D:354:ALA:N	2.30	0.43
2:D:105:ALA:O	2:D:109:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:VAL:O	1:A:68:GLY:O	2.37	0.43
1:B:107:VAL:HG23	1:B:202:ILE:CD1	2.48	0.42
1:A:330:ALA:HA	1:A:340:ILE:HG21	2.01	0.42
1:B:233:ALA:HB2	2:D:145:ALA:CB	2.49	0.42
1:B:224:LYS:HB2	2:D:142:LEU:O	2.18	0.42
1:B:451:HIS:HB3	1:B:463:GLU:HB2	2.01	0.42
1:A:304:ALA:HB1	1:A:483:LYS:NZ	2.34	0.42
2:C:215:ARG:HB3	2:C:218:THR:HG1	1.84	0.42
1:B:18:VAL:HG12	1:B:55:GLY:O	2.18	0.42
1:A:89:ASN:ND2	5:A:1003:N8E:H132	2.34	0.42
2:D:30:ARG:NH1	2:D:67:GLU:HG2	2.35	0.42
1:A:226:ASN:OD1	1:A:227:ALA:N	2.51	0.42
2:C:324:LYS:NZ	2:C:324:LYS:HB3	2.34	0.42
1:B:1:MET:SD	1:B:4:GLU:HB2	2.59	0.42
1:A:404:ASN:HB3	1:A:407:VAL:HB	2.01	0.42
1:B:507:TYR:CE1	1:B:631:MET:HB3	2.54	0.42
1:B:628:ILE:O	1:B:632:MET:HG2	2.20	0.42
1:A:408:LYS:HD2	1:A:432:ILE:HD12	2.02	0.42
1:B:178:ASP:O	1:B:182:VAL:HG22	2.19	0.42
2:C:189:PHE:HB3	2:C:193:ILE:CD1	2.49	0.42
2:D:31:ALA:HB2	2:D:124:HIS:HB2	2.01	0.42
1:B:56:VAL:O	1:B:106:THR:HA	2.20	0.42
1:B:32:PHE:CE1	1:B:66:ILE:HG21	2.55	0.42
1:A:638:GLU:HA	1:A:641:ARG:NH1	2.35	0.42
2:C:357:SER:O	2:C:360:LEU:HB3	2.19	0.42
2:D:49:VAL:HG23	2:D:49:VAL:O	2.20	0.42
2:C:286:TYR:CZ	2:C:289:VAL:HG11	2.55	0.42
2:D:292:THR:O	2:D:296:LEU:HG	2.19	0.42
1:A:358:ALA:HB1	1:B:376:ALA:HB1	2.00	0.42
1:A:475:VAL:O	1:A:479:VAL:HG23	2.18	0.42
1:A:111:ASN:HA	1:A:194:LEU:CD1	2.49	0.42
1:A:226:ASN:HB3	1:A:229:GLU:H	1.84	0.42
1:B:499:LEU:C	1:B:499:LEU:HD13	2.40	0.42
1:A:628:ILE:O	1:A:632:MET:HG2	2.20	0.42
1:B:77:ASN:O	1:B:85:LEU:HD13	2.18	0.42
1:A:401:VAL:HG11	1:A:407:VAL:HG12	2.01	0.42
1:B:434:ILE:HG23	1:B:447:PHE:CE2	2.55	0.42
1:A:104:VAL:HG22	1:A:105:PRO:HD2	2.02	0.42
1:A:499:LEU:C	1:A:499:LEU:HD13	2.40	0.42
1:A:176:ALA:HB1	1:A:188:VAL:HG11	2.01	0.42
2:D:269:ALA:HB2	2:D:361:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:296:LEU:HD21	2:D:306:ILE:HD11	2.02	0.42
1:B:325:MET:O	1:B:329:ILE:HG13	2.19	0.42
2:C:299:ALA:O	2:C:301:LEU:HG	2.20	0.42
1:A:546:ILE:N	1:A:546:ILE:HD12	2.35	0.42
1:A:702:LYS:O	1:A:706:MET:HG2	2.20	0.42
1:A:464:VAL:HG21	1:A:479:VAL:HG22	2.01	0.42
1:B:700:THR:OG1	1:B:703:LEU:HB2	2.20	0.42
1:A:658:LEU:HB3	1:A:662:ILE:HD11	2.01	0.41
2:D:43:LEU:HD11	2:D:82:ILE:HD11	2.01	0.41
2:D:6:ARG:HB3	2:D:110:MET:HA	2.02	0.41
1:A:157:ARG:HG2	1:A:223:LEU:HD23	2.02	0.41
2:C:39:ILE:HG23	2:C:118:VAL:HG11	2.02	0.41
1:B:509:GLY:O	1:B:513:LYS:HG3	2.20	0.41
1:B:540:LEU:HD23	1:B:544:VAL:HG23	2.02	0.41
1:A:57:ILE:HG12	1:A:202:ILE:HG13	2.02	0.41
2:D:313:GLU:OE2	2:D:341:GLY:HA3	2.20	0.41
1:B:655:ASP:HB3	1:B:670:GLY:HA3	2.02	0.41
1:A:350:ILE:HD11	1:A:386:LEU:HG	2.01	0.41
2:C:125:MET:CE	2:C:246:THR:H	2.34	0.41
1:A:146:TYR:HB2	1:A:147:PRO:HD2	2.01	0.41
1:B:157:ARG:NH1	1:B:223:LEU:HG	2.35	0.41
1:A:18:VAL:HG12	1:A:55:GLY:O	2.20	0.41
2:D:329:LEU:O	2:D:329:LEU:HD23	2.20	0.41
1:B:672:LEU:HA	1:B:675:ILE:HD12	2.02	0.41
2:C:307:ASP:O	2:C:335:LYS:HB2	2.20	0.41
2:D:273:SER:OG	2:D:299:ALA:HB2	2.21	0.41
1:A:525:ASP:OD2	1:A:538:ALA:HB3	2.20	0.41
1:B:464:VAL:HG11	1:B:475:VAL:HG13	2.03	0.41
1:B:185:VAL:HG12	1:B:187:ALA:H	1.85	0.41
1:A:93:ASN:OD1	1:A:149:PHE:HA	2.19	0.41
2:C:293:GLN:O	2:C:297:LYS:HG3	2.20	0.41
1:A:141:VAL:HG12	1:A:251:ALA:CB	2.46	0.41
2:D:172:PHE:CE2	2:D:313:GLU:HG2	2.54	0.41
1:A:56:VAL:O	1:A:106:THR:HA	2.21	0.41
2:C:357:SER:O	2:C:361:LEU:HG	2.21	0.41
1:A:384:PRO:HB2	1:B:383:ARG:HD3	2.02	0.41
1:B:127:PHE:CD2	1:B:201:LEU:HD21	2.56	0.41
1:A:338:THR:OG1	1:A:484:LYS:HE3	2.21	0.41
1:B:157:ARG:HD2	1:B:219:LYS:HA	2.02	0.41
1:A:461:LEU:HB2	1:A:660:TYR:CB	2.50	0.41
1:B:525:ASP:O	1:B:529:GLU:HG3	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:ASP:CG	1:B:567:ASP:H	2.24	0.41
1:A:344:ASP:OD1	1:A:345:ILE:N	2.52	0.41
1:A:532:GLY:C	1:A:665:PRO:HG3	2.40	0.41
1:B:681:ALA:HB3	1:B:682:GLU:OE1	2.21	0.41
2:D:158:LEU:HD23	2:D:318:GLN:HA	2.03	0.41
2:C:93:ARG:O	2:C:93:ARG:HG2	2.21	0.41
2:C:199:TYR:HA	2:C:204:PHE:O	2.21	0.41
1:B:371:THR:HG22	1:B:373:ALA:H	1.85	0.41
1:B:228:ILE:HG22	2:D:282:ALA:CB	2.50	0.41
2:D:164:ILE:CG2	2:D:169:GLN:HG3	2.51	0.41
1:B:408:LYS:HD2	1:B:432:ILE:HD12	2.02	0.41
2:D:181:HIS:O	2:D:185:VAL:HG23	2.20	0.41
1:A:320:LEU:HB2	1:A:399:GLU:HA	2.02	0.41
2:D:347:HIS:CD2	3:D:4001:ACO:S1P	3.14	0.41
1:A:70:ASP:HA	1:A:294:PHE:HZ	1.86	0.41
1:B:78:PHE:HA	1:B:85:LEU:HD13	2.03	0.41
1:B:403:GLU:HA	1:B:408:LYS:HE2	2.03	0.41
2:C:206:LYS:HB2	2:C:206:LYS:NZ	2.36	0.41
1:A:404:ASN:ND2	5:A:1004:N8E:H031	2.31	0.40
2:D:58:ILE:HD13	2:D:104:THR:HB	2.02	0.40
2:C:98:SER:HB3	2:C:354:ALA:H	1.86	0.40
2:C:61:CYS:SG	2:C:68:GLN:HB3	2.60	0.40
2:D:155:ALA:HB2	2:D:315:PHE:CD2	2.57	0.40
1:B:539:TYR:O	1:B:543:VAL:HG23	2.21	0.40
1:A:386:LEU:HB2	1:B:390:ASP:OD1	2.20	0.40
1:B:445:GLU:HA	1:B:469:LYS:O	2.21	0.40
2:C:121:GLY:HA3	2:C:350:GLY:O	2.21	0.40
2:C:122:VAL:HG22	2:C:123:GLU:N	2.35	0.40
1:A:140:GLU:OE2	5:A:1003:N8E:H051	2.21	0.40
1:A:105:PRO:HG3	1:A:212:TYR:CD2	2.57	0.40
1:A:443:ARG:HA	1:A:445:GLU:OE2	2.21	0.40
1:A:388:TYR:CE2	1:A:415:VAL:HG22	2.55	0.40
1:A:127:PHE:CD2	1:A:201:LEU:HD21	2.56	0.40
1:B:226:ASN:O	1:B:230:GLN:HG3	2.20	0.40
2:D:74:ARG:CZ	2:D:78:LEU:HD21	2.52	0.40
2:C:189:PHE:HB3	2:C:193:ILE:HD12	2.03	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	713/715 (100%)	652 (91%)	52 (7%)	9 (1%)	15	60
1	B	702/715 (98%)	637 (91%)	56 (8%)	9 (1%)	15	60
2	C	388/390 (100%)	353 (91%)	32 (8%)	3 (1%)	24	70
2	D	388/390 (100%)	350 (90%)	36 (9%)	2 (0%)	34	78
All	All	2191/2210 (99%)	1992 (91%)	176 (8%)	23 (1%)	19	66

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	GLY
1	A	250	PRO
1	B	250	PRO
1	B	286	ALA
1	B	68	GLY
1	B	566	ASP
1	B	567	ASP
2	C	349	PHE
1	A	289	CYS
1	A	597	LYS
2	C	332	MET
2	D	349	PHE
1	A	65	PHE
1	A	252	PRO
1	B	285	SER
2	D	332	MET
1	A	500	VAL
1	B	324	ILE
1	B	496	PRO
2	C	54	VAL
1	A	324	ILE
1	A	161	VAL
1	B	252	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	547/562 (97%)	526 (96%)	21 (4%)	40	76
1	B	530/562 (94%)	516 (97%)	14 (3%)	54	83
2	C	305/307 (99%)	289 (95%)	16 (5%)	29	68
2	D	304/307 (99%)	294 (97%)	10 (3%)	45	79
All	All	1686/1738 (97%)	1625 (96%)	61 (4%)	42	77

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LEU
1	A	89	ASN
1	A	104	VAL
1	A	140	GLU
1	A	158	LEU
1	A	194	LEU
1	A	199	LEU
1	A	217	GLN
1	A	247	PRO
1	A	248	ASN
1	A	249	TYR
1	A	280	LYS
1	A	296	ASN
1	A	298	GLN
1	A	366	ASP
1	A	421	GLU
1	A	445	GLU
1	A	525	ASP
1	A	540	LEU
1	A	554	ASP
1	A	676	ASP
1	B	38	ASN
1	B	40	LEU
1	B	74	PHE

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Mol	Chain	Res	Type
1	B	104	VAL
1	B	119	LEU
1	B	194	LEU
1	B	290	LEU
1	B	295	LEU
1	B	366	ASP
1	B	390	ASP
1	B	473	LEU
1	B	525	ASP
1	B	643	LEU
1	B	676	ASP
2	C	3	LEU
2	C	4	ASN
2	C	5	PRO
2	C	7	ASP
2	C	27	ARG
2	C	150	MET
2	C	201	GLU
2	C	206	LYS
2	C	231	ASN
2	C	265	LEU
2	C	284	MET
2	C	289	VAL
2	C	318	GLN
2	C	325	ASP
2	C	351	CYS
2	C	355	ARG
2	D	55	GLU
2	D	175	ARG
2	D	284	MET
2	D	289	VAL
2	D	318	GLN
2	D	325	ASP
2	D	330	ASP
2	D	339	HIS
2	D	351	CYS
2	D	355	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN

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Mol	Chain	Res	Type
1	A	42	GLN
1	A	77	ASN
1	A	89	ASN
1	A	288	ASN
1	A	296	ASN
1	A	298	GLN
1	A	404	ASN
1	A	446	ASN
1	A	501	ASN
1	A	629	ASN
1	A	698	HIS
1	B	38	ASN
1	B	42	GLN
1	B	296	ASN
1	B	488	ASN
1	B	629	ASN
2	C	4	ASN
2	C	37	HIS
2	C	46	ASN
2	C	113	ASN
2	C	162	HIS
2	C	169	GLN
2	C	197	GLN
2	C	202	ASN
2	C	231	ASN
2	C	362	ASN
2	D	4	ASN
2	D	46	ASN
2	D	162	HIS
2	D	197	GLN
2	D	293	GLN
2	D	362	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	NAD	A	1001	-	31,38,48	0.76	0	39,58,73	1.03	3 (7%)
5	N8E	A	1002	-	23,23,23	0.43	0	22,22,22	0.43	0
5	N8E	A	1003	-	23,23,23	0.41	0	22,22,22	0.42	0
5	N8E	A	1004	-	23,23,23	0.45	0	22,22,22	0.42	0
4	NAD	B	2001	-	38,48,48	1.31	3 (7%)	47,73,73	1.78	8 (17%)
3	ACO	C	3001	-	32,42,53	0.83	0	40,66,79	1.02	3 (7%)
3	ACO	D	4001	-	43,53,53	0.84	0	55,79,79	0.95	4 (7%)

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	NAD	A	1001	-	-	0/18/51/62	0/4/4/5
5	N8E	A	1002	-	-	0/21/21/21	0/0/0/0
5	N8E	A	1003	-	-	0/21/21/21	0/0/0/0
5	N8E	A	1004	-	-	0/21/21/21	0/0/0/0
4	NAD	B	2001	-	-	0/22/62/62	0/5/5/5
3	ACO	C	3001	-	-	0/33/54/67	0/3/3/3
3	ACO	D	4001	-	-	2/47/67/67	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2001	NAD	C6N-N1N	2.94	1.43	1.35
4	B	2001	NAD	C4N-C3N	3.02	1.44	1.39
4	B	2001	NAD	C3N-C7N	5.08	1.58	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2001	NAD	C5N-C4N-C3N	-5.29	113.69	120.33
3	C	3001	ACO	P2A-O3A-P1A	-4.28	120.72	132.73
3	D	4001	ACO	P2A-O3A-P1A	-4.19	120.97	132.73
4	A	1001	NAD	PN-O3-PA	-4.16	121.04	132.73
4	B	2001	NAD	PN-O3-PA	-4.11	121.19	132.73
4	B	2001	NAD	C5N-C6N-N1N	-3.17	114.99	120.47
3	D	4001	ACO	C2P-S1P-C	2.01	114.27	101.83
3	C	3001	ACO	O3A-P2A-O6A	2.31	109.08	102.94
3	D	4001	ACO	O3A-P1A-O5B	2.32	109.10	102.94
4	B	2001	NAD	O3-PN-O5D	2.38	109.25	102.94
4	A	1001	NAD	O3-PN-O5D	2.39	109.29	102.94
3	C	3001	ACO	O3A-P1A-O5B	2.40	109.30	102.94
4	A	1001	NAD	O3-PA-O5B	2.42	109.34	102.94
4	B	2001	NAD	O3-PA-O5B	2.42	109.36	102.94
3	D	4001	ACO	O3A-P2A-O6A	2.42	109.36	102.94
4	B	2001	NAD	C2N-C3N-C4N	3.13	121.77	118.29
4	B	2001	NAD	O4D-C1D-N1N	5.19	113.83	108.13
4	B	2001	NAD	C6N-C5N-C4N	5.34	127.52	119.44

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	4001	ACO	CH3-C-S1P-C2P
3	D	4001	ACO	O-C-S1P-C2P

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1002	N8E	1	0
5	A	1003	N8E	2	0
5	A	1004	N8E	3	0
4	B	2001	NAD	1	0
3	D	4001	ACO	4	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	715/715 (100%)	-0.37	9 (1%) 79 70	1, 8, 48, 139	0
1	B	706/715 (98%)	-0.15	11 (1%) 74 65	1, 25, 71, 130	0
2	C	390/390 (100%)	-0.51	0 100 100	1, 4, 27, 70	0
2	D	390/390 (100%)	-0.52	1 (0%) 94 91	1, 3, 30, 91	0
All	All	2201/2210 (99%)	-0.35	21 (0%) 84 76	1, 9, 58, 139	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	598	GLY	6.7
1	A	248	ASN	4.1
2	D	2	SER	3.9
1	A	597	LYS	3.7
1	A	366	ASP	3.4
1	A	595	ASP	3.4
1	B	606	SER	3.0
1	B	715	GLY	2.7
1	B	285	SER	2.6
1	B	287	SER	2.6
1	B	562	ASP	2.6
1	B	710	GLY	2.6
1	B	247	PRO	2.5
1	A	367	LYS	2.3
1	B	368	GLY	2.3
1	A	348	HIS	2.3
1	B	295	LEU	2.2
1	B	566	ASP	2.2
1	B	243	GLY	2.2
1	A	368	GLY	2.1
1	A	715	GLY	2.1

## 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( $\text{\AA}^2$ )	Q<0.9
5	N8E	A	1002	24/24	0.74	0.43	5.06	23,23,23,23	0
3	ACO	C	3001	40/51	0.70	0.40	4.36	67,80,128,128	0
3	ACO	D	4001	51/51	0.77	0.39	4.34	62,68,133,133	0
5	N8E	A	1004	24/24	0.62	0.47	4.12	26,58,77,77	0
5	N8E	A	1003	24/24	0.82	0.29	1.59	15,15,15,15	0
4	NAD	B	2001	44/44	0.80	0.36	1.41	43,56,93,93	0
4	NAD	A	1001	35/44	0.83	0.31	0.97	20,44,116,116	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.