



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 01:21 PM GMT

PDB ID : 2IOA  
Title : E. coli Bifunctional glutathionylspermidinesynthetase/amidase Incomplex with Mg<sup>2+</sup> and ADP and phosphinate inhibitor  
Authors : Pai, C.H.; Chiang, B.Y.; Ko, T.P.; Chou, C.C.; Chong, C.M.; Yen, F.J.; Coward, J.K.; Wang, A.H.-J.; Lin, C.H.  
Deposited on : 2006-10-10  
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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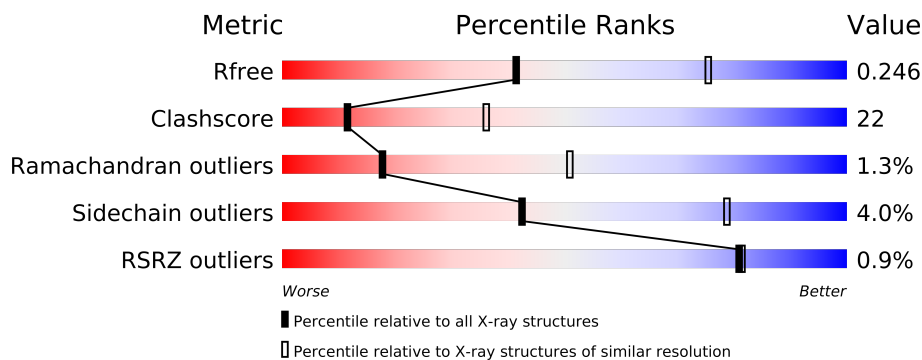
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	619	
1	B	619	

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 10089 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

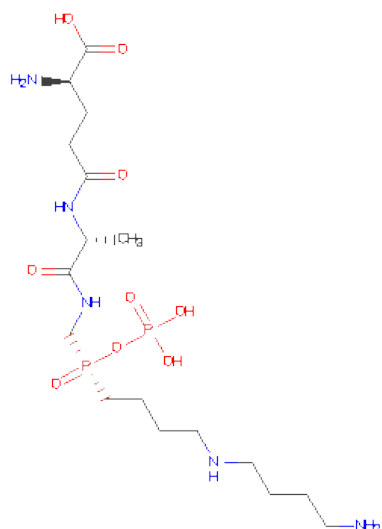
- Molecule 1 is a protein called Bifunctional glutathionylspermidinesynthetase/amidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	589	Total	C	N	O	S	0	0	0
			4755	3051	813	872	19			
1	B	589	Total	C	N	O	S	0	0	0
			4753	3048	813	873	19			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

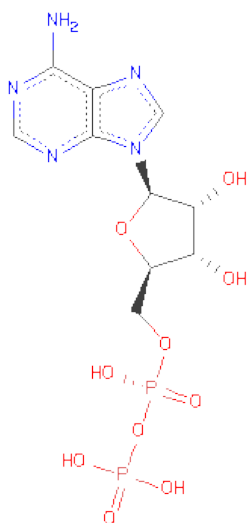
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is D-GAMMA-GLUTAMYL-N-{[(R)-{4-[(4-AMINO BUTYL)AMINO] BUTYL } (PHOSPHONOOXY)PHOSPHORYL] METHYL}-D-ALANINAMIDE (three-letter code: GGA) (formula: C<sub>17</sub>H<sub>37</sub>N<sub>5</sub>O<sub>9</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			33	17	5	9	2		
3	B	1	Total	C	N	O	P	0	0
			33	17	5	9	2		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is water.

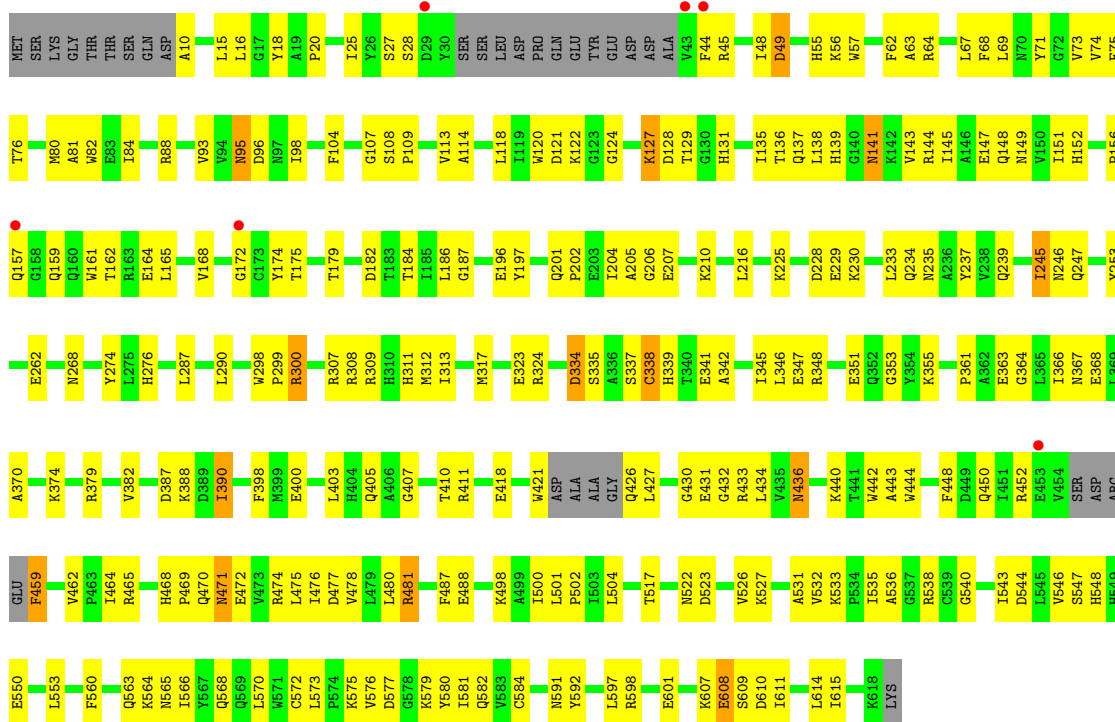
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	236	Total	O	0	0
			236	236		
5	B	221	Total	O	0	0
			221	221		

### 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.

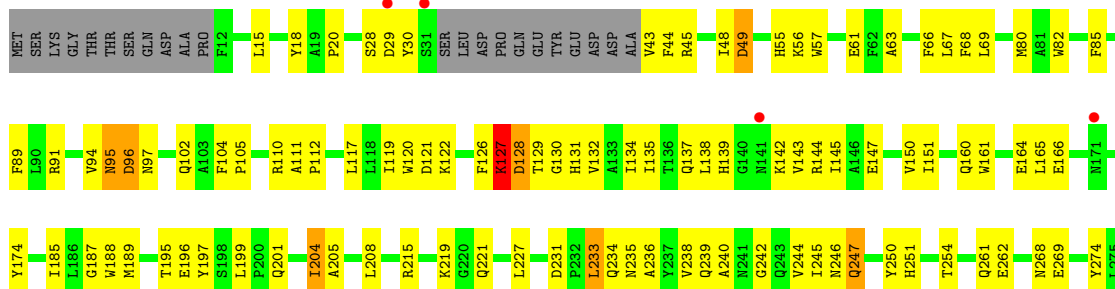
- Molecule 1: Bifunctional glutathionylspermidinesynthetase/amidase

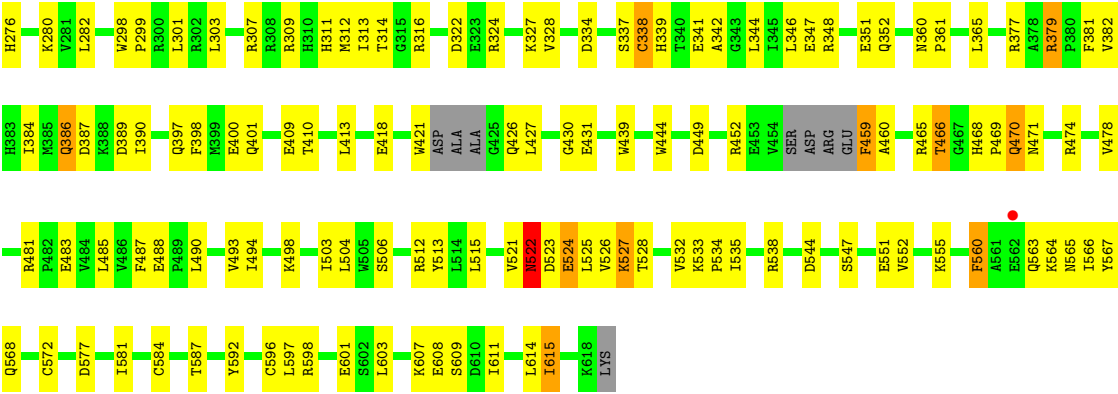
Chain A: 



- Molecule 1: Bifunctional glutathionylspermidinesynthetase/amidase

Chain B: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.03Å 75.29Å 84.66Å 70.09° 74.06° 77.55°	Depositor
Resolution (Å)	30.00 – 2.80 29.49 – 2.69	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.80) 93.6 (29.49-2.69)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.23 (at 2.68Å)	Xtriage
Refinement program	XTALVIEW	Depositor
R, $R_{free}$	0.173 , 0.243 0.184 , 0.246	Depositor DCC
$R_{free}$ test set	3420 reflections (11.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.8	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 32.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 35215 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10089	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, GGA, ADP

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	0.43	0/4877	0.67	0/6621
1	B	0.43	0/4874	0.68	0/6615
All	All	0.43	0/9751	0.68	0/13236

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4755	0	4641	217	0
1	B	4753	0	4637	202	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	33	0	35	8	0
3	B	33	0	34	2	0
4	A	27	0	11	2	0
4	B	27	0	11	0	0
5	A	236	0	0	18	0
5	B	221	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10089	0	9369	417	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (417) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:245:ILE:O	5:A:4192:HOH:O	1.73	1.03
1:A:210:LYS:HD3	1:A:323:GLU:HB3	1.40	1.02
1:A:64:ARG:HG3	1:A:74:VAL:HG23	1.43	0.96
1:A:311:HIS:ND1	5:A:4195:HOH:O	1.97	0.96
1:A:341:GLU:HG2	1:A:611:ILE:HG13	1.48	0.93
1:B:608:GLU:CD	1:B:608:GLU:H	1.79	0.87
1:B:268:ASN:HD21	1:B:592:TYR:H	1.24	0.83
1:A:311:HIS:CE1	5:A:4195:HOH:O	2.32	0.83
1:B:607:LYS:O	3:B:6002:GGA:H432	1.79	0.83
1:B:465:ARG:HG3	1:B:474:ARG:HH21	1.42	0.82
1:B:233:LEU:HD12	1:B:401:GLN:OE1	1.80	0.81
1:A:607:LYS:O	3:A:6001:GGA:H432	1.82	0.80
1:B:337:SER:O	1:B:338:CYS:HB2	1.81	0.80
1:A:121:ASP:HB2	1:A:186:LEU:HD21	1.62	0.80
1:A:28:SER:HB3	1:A:57:TRP:HB2	1.63	0.79
1:B:341:GLU:HG2	1:B:611:ILE:HG13	1.65	0.79
1:B:384:ILE:HG23	1:B:439:TRP:HE3	1.49	0.78
1:B:418:GLU:HG3	1:B:430:GLY:H	1.49	0.77
1:A:311:HIS:O	1:A:313:ILE:HD12	1.86	0.75
1:A:442:TRP:HE3	3:A:6001:GGA:O2	1.67	0.75
1:A:498:LYS:HB3	1:A:566:ILE:HD13	1.70	0.74
1:A:390:ILE:H	1:A:390:ILE:HD12	1.51	0.74
1:B:245:ILE:HG13	1:B:246:ASN:H	1.52	0.74
1:A:540:GLY:HA2	1:A:543:ILE:HD12	1.71	0.72
1:B:361:PRO:HG3	1:B:614:LEU:HD23	1.71	0.72
1:B:138:LEU:HD12	1:B:143:VAL:HG12	1.71	0.72
1:A:337:SER:O	1:A:338:CYS:HB2	1.88	0.72
1:A:138:LEU:HD12	1:A:143:VAL:HG12	1.71	0.72
1:A:137:GLN:HG2	1:A:144:ARG:HB2	1.72	0.71
1:A:421:TRP:H	1:A:471:ASN:HD21	1.38	0.71
1:A:442:TRP:CE3	3:A:6001:GGA:O2	2.44	0.71
1:B:204:ILE:HD13	1:B:205:ALA:N	2.05	0.71
1:A:151:ILE:HD12	1:A:162:THR:HB	1.72	0.70
1:B:301:LEU:HD13	1:B:494:ILE:HD11	1.73	0.70
1:A:48:ILE:HD13	1:A:69:LEU:HD21	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:96:ASP:CG	1:A:309:ARG:HH21	1.96	0.69
1:A:137:GLN:HE21	1:A:139:HIS:CE1	2.10	0.69
1:A:268:ASN:HD21	1:A:592:TYR:H	1.40	0.69
1:A:448:PHE:O	1:A:452:ARG:HG3	1.93	0.69
1:A:20:PRO:HG2	1:A:136:THR:HB	1.74	0.69
1:A:579:LYS:HD2	5:A:4086:HOH:O	1.92	0.68
1:A:201:GLN:HG3	1:A:202:PRO:HD2	1.73	0.68
1:A:157:GLN:CD	1:A:157:GLN:H	1.96	0.68
1:B:413:LEU:HD23	1:B:418:GLU:OE1	1.94	0.68
1:A:82:TRP:HD1	1:A:129:THR:O	1.77	0.68
1:B:532:VAL:HB	1:B:544:ASP:HB2	1.74	0.68
1:B:245:ILE:HG13	1:B:246:ASN:N	2.08	0.67
1:A:379:ARG:HH12	1:A:436:ASN:HB2	1.60	0.67
1:A:575:LYS:HD2	1:A:580:TYR:CZ	2.29	0.67
1:B:522:ASN:HD22	1:B:522:ASN:C	1.98	0.67
1:B:122:LYS:HB2	1:B:127:LYS:O	1.95	0.67
1:A:337:SER:OG	3:A:6001:GGA:H372	1.95	0.66
1:A:418:GLU:HG2	1:A:418:GLU:O	1.94	0.66
1:A:527:LYS:O	1:A:548:HIS:HB3	1.95	0.66
1:B:233:LEU:HD11	1:B:398:PHE:HA	1.79	0.65
1:B:418:GLU:CG	1:B:430:GLY:H	2.09	0.65
1:B:598:ARG:NH1	5:B:4308:HOH:O	2.28	0.65
1:A:535:ILE:HG22	1:A:566:ILE:HG23	1.79	0.65
1:A:540:GLY:HA2	1:A:543:ILE:CD1	2.27	0.65
1:A:135:ILE:HD13	1:A:145:ILE:HG22	1.78	0.65
1:B:522:ASN:HD21	1:B:525:LEU:H	1.43	0.65
1:A:341:GLU:CG	1:A:611:ILE:HG13	2.24	0.64
1:A:210:LYS:HD3	1:A:323:GLU:CB	2.22	0.64
1:A:95:ASN:HD22	1:A:95:ASN:N	1.94	0.64
1:B:227:LEU:HD23	1:B:348:ARG:HG2	1.78	0.64
1:A:64:ARG:CG	1:A:74:VAL:HG23	2.21	0.64
1:A:426:GLN:N	5:A:4229:HOH:O	2.30	0.63
1:A:15:LEU:HD21	1:A:18:TYR:CZ	2.33	0.63
1:B:215:ARG:HD3	5:B:4401:HOH:O	1.98	0.63
1:B:102:GLN:NE2	1:B:195:THR:HG22	2.14	0.63
1:B:384:ILE:HG23	1:B:439:TRP:CE3	2.34	0.63
1:B:303:LEU:HD13	1:B:307:ARG:NH2	2.14	0.63
1:A:411:ARG:NE	1:A:418:GLU:OE1	2.28	0.62
1:A:579:LYS:HE3	1:A:601:GLU:OE2	1.98	0.62
1:B:164:GLU:O	1:B:165:LEU:HD23	1.99	0.62
1:B:105:PRO:HD3	1:B:199:LEU:HG	1.79	0.62
1:A:210:LYS:NZ	1:A:323:GLU:HG2	2.13	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:298:TRP:HB2	1:B:299:PRO:HD3	1.81	0.62
1:B:137:GLN:CD	1:B:144:ARG:HD2	2.20	0.62
1:A:390:ILE:N	1:A:390:ILE:HD12	2.15	0.61
1:B:498:LYS:HB2	1:B:535:ILE:HA	1.80	0.61
1:A:124:GLY:HA3	1:A:182:ASP:O	2.00	0.61
1:B:418:GLU:HG3	1:B:430:GLY:N	2.14	0.61
1:A:144:ARG:HD3	1:A:161:TRP:CD1	2.35	0.61
1:B:204:ILE:HD13	1:B:205:ALA:H	1.63	0.61
1:A:462:VAL:HG12	1:A:464:ILE:HG23	1.81	0.61
1:A:400:GLU:HG3	1:A:410:THR:OG1	2.01	0.60
1:B:166:GLU:HB3	5:B:4283:HOH:O	2.00	0.60
1:A:121:ASP:HB2	1:A:186:LEU:CD2	2.32	0.60
1:A:390:ILE:H	1:A:390:ILE:CD1	2.12	0.60
1:B:246:ASN:HB2	5:B:4468:HOH:O	2.02	0.60
1:B:498:LYS:HB3	1:B:566:ILE:HD13	1.84	0.60
1:A:474:ARG:NH2	1:A:477:ASP:OD2	2.35	0.60
1:A:298:TRP:HB2	1:A:299:PRO:HD3	1.84	0.60
1:B:135:ILE:HA	1:B:145:ILE:HG22	1.82	0.60
1:B:465:ARG:HG3	1:B:474:ARG:NH2	2.16	0.59
1:B:522:ASN:O	1:B:526:VAL:HG23	2.02	0.59
1:A:137:GLN:CD	1:A:144:ARG:HD2	2.23	0.59
1:B:247:GLN:HG3	5:B:4468:HOH:O	2.03	0.59
1:A:120:TRP:HB2	1:A:131:HIS:HB3	1.84	0.58
1:A:341:GLU:HG2	1:A:611:ILE:CG1	2.29	0.58
1:A:104:PHE:O	1:A:187:GLY:HA3	2.02	0.58
1:A:573:LEU:HD22	1:A:581:ILE:O	2.04	0.58
1:B:382:VAL:HB	1:B:410:THR:HG22	1.84	0.58
1:B:238:VAL:O	1:B:242:GLY:N	2.32	0.58
1:B:82:TRP:HD1	1:B:129:THR:O	1.86	0.58
1:B:523:ASP:O	1:B:526:VAL:HB	2.04	0.58
1:B:102:GLN:CD	1:B:195:THR:HG22	2.24	0.58
1:B:251:HIS:HE1	1:B:577:ASP:OD2	1.86	0.58
1:A:95:ASN:N	1:A:95:ASN:ND2	2.52	0.57
1:B:48:ILE:O	1:B:49:ASP:HB2	2.04	0.57
1:A:137:GLN:HE21	1:A:139:HIS:HE1	1.50	0.57
1:A:387:ASP:OD1	1:A:388:LYS:N	2.36	0.57
1:A:235:ASN:O	1:A:239:GLN:HG2	2.04	0.57
1:A:361:PRO:HG3	1:A:614:LEU:HD23	1.86	0.57
1:A:121:ASP:CB	1:A:186:LEU:HD21	2.33	0.57
1:B:538:ARG:NH1	1:B:560:PHE:HZ	2.03	0.57
1:B:235:ASN:O	1:B:239:GLN:HG2	2.05	0.57
1:B:341:GLU:CG	1:B:611:ILE:HG13	2.34	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:403:LEU:HD13	1:A:410:THR:HG21	1.87	0.56
1:A:312:MET:HA	1:A:488:GLU:OE2	2.06	0.56
1:A:351:GLU:O	1:A:351:GLU:HG2	2.05	0.56
1:B:282:LEU:HD21	1:B:494:ILE:CD1	2.35	0.56
1:B:468:HIS:CG	1:B:469:PRO:HD2	2.40	0.56
1:A:538:ARG:NH1	1:A:560:PHE:HZ	2.02	0.56
1:B:94:VAL:HG13	5:B:4347:HOH:O	2.05	0.56
1:B:322:ASP:OD2	1:B:324:ARG:HB2	2.05	0.56
1:A:164:GLU:O	1:A:165:LEU:HD23	2.05	0.56
1:B:522:ASN:ND2	1:B:525:LEU:H	2.02	0.56
1:B:303:LEU:HD13	1:B:307:ARG:HH21	1.70	0.56
1:A:502:PRO:HG3	1:A:517:THR:HG22	1.88	0.56
1:A:131:HIS:HE1	1:A:147:GLU:OE2	1.89	0.55
1:A:347:GLU:OE2	1:A:363:GLU:HB3	2.06	0.55
1:A:335:SER:HB2	5:A:4005:HOH:O	2.05	0.55
1:B:44:PHE:HD2	1:B:55:HIS:HE1	1.54	0.55
1:A:151:ILE:CD1	1:A:162:THR:HB	2.35	0.55
1:A:225:LYS:HE2	1:A:229:GLU:OE2	2.07	0.55
1:A:80:MET:SD	1:A:129:THR:HG22	2.46	0.55
1:B:127:LYS:HB3	1:B:128:ASP:OD1	2.07	0.55
1:B:112:PRO:HB3	1:B:188:TRP:CD2	2.41	0.55
1:A:307:ARG:HD3	1:B:49:ASP:OD2	2.07	0.55
1:B:63:ALA:O	1:B:67:LEU:HG	2.07	0.55
1:A:465:ARG:NH2	1:A:472:GLU:O	2.38	0.55
1:A:107:GLY:HA2	1:A:174:TYR:O	2.07	0.55
1:A:137:GLN:NE2	1:A:144:ARG:HD2	2.22	0.54
1:A:426:GLN:HG3	5:A:4381:HOH:O	2.06	0.54
1:A:597:LEU:HD11	1:A:614:LEU:HD13	1.89	0.54
1:A:44:PHE:HB3	1:A:55:HIS:CE1	2.42	0.54
1:A:114:ALA:HB1	1:B:460:ALA:HB2	1.88	0.54
1:B:481:ARG:NH1	1:B:483:GLU:OE1	2.39	0.54
1:B:572:CYS:HB3	1:B:603:LEU:HD22	1.90	0.54
1:A:113:VAL:HG22	1:A:114:ALA:N	2.22	0.54
1:B:80:MET:SD	1:B:129:THR:HG22	2.47	0.54
1:B:301:LEU:HD23	1:B:490:LEU:HG	1.89	0.54
1:B:89:PHE:CE2	1:B:269:GLU:HG3	2.42	0.54
1:A:345:ILE:N	1:A:345:ILE:HD12	2.23	0.54
1:A:427:LEU:HD21	1:A:478:VAL:HG22	1.90	0.54
1:A:335:SER:OG	3:A:6001:GGA:H223	2.07	0.54
1:A:48:ILE:O	1:A:49:ASP:HB2	2.08	0.54
1:A:433:ARG:HG2	1:A:434:LEU:N	2.23	0.54
1:A:71:TYR:HB2	1:A:73:VAL:HG22	1.89	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:546:VAL:O	1:A:570:LEU:HD22	2.08	0.53
1:A:540:GLY:CA	1:A:543:ILE:HD12	2.36	0.53
1:A:450:GLN:OE1	1:A:475:LEU:HB3	2.09	0.53
1:A:96:ASP:OD1	1:A:309:ARG:NH2	2.38	0.53
1:B:45:ARG:NH2	5:B:4203:HOH:O	2.33	0.53
1:B:254:THR:HA	1:B:615:ILE:O	2.08	0.53
1:B:311:HIS:N	1:B:311:HIS:CD2	2.76	0.53
1:A:532:VAL:HB	1:A:544:ASP:HB2	1.89	0.52
1:A:342:ALA:HA	1:A:346:LEU:HD12	1.91	0.52
1:A:207:GLU:OE2	1:A:207:GLU:HA	2.09	0.52
1:B:268:ASN:ND2	1:B:592:TYR:H	2.01	0.52
1:A:45:ARG:O	1:A:56:LYS:HE3	2.10	0.52
1:A:96:ASP:OD1	1:A:309:ARG:NE	2.40	0.52
1:B:44:PHE:HD2	1:B:55:HIS:CE1	2.26	0.52
1:A:81:ALA:O	1:A:84:ILE:HG13	2.09	0.52
1:B:533:LYS:NZ	1:B:568:GLN:HE22	2.08	0.52
1:B:236:ALA:CB	1:B:397:GLN:HE22	2.22	0.52
1:A:122:LYS:HB2	1:A:127:LYS:O	2.09	0.52
1:B:527:LYS:HE2	1:B:528:THR:HG22	1.92	0.52
1:A:98:ILE:HD12	1:A:98:ILE:N	2.25	0.52
1:A:95:ASN:HD22	1:A:95:ASN:H	1.56	0.52
1:A:196:GLU:O	1:A:197:TYR:HB2	2.10	0.52
1:B:339:HIS:H	1:B:339:HIS:CD2	2.27	0.52
1:A:57:TRP:HB3	1:A:148:GLN:NE2	2.24	0.52
1:B:254:THR:HG22	1:B:615:ILE:HG22	1.91	0.51
1:A:210:LYS:HZ2	1:A:323:GLU:HG2	1.73	0.51
1:B:57:TRP:HZ3	1:B:61:GLU:OE1	1.93	0.51
1:B:342:ALA:HA	1:B:346:LEU:HD12	1.92	0.51
1:B:139:HIS:HB2	1:B:142:LYS:O	2.10	0.51
1:B:341:GLU:CB	1:B:611:ILE:HG13	2.41	0.51
1:B:126:PHE:O	1:B:127:LYS:O	2.28	0.51
1:B:144:ARG:HD3	1:B:161:TRP:CD1	2.45	0.51
1:B:534:PRO:HB2	1:B:563:GLN:NE2	2.26	0.51
1:A:370:ALA:O	1:A:374:LYS:HG3	2.11	0.51
1:B:328:VAL:HG12	1:B:513:TYR:O	2.11	0.51
1:A:118:LEU:HD22	1:A:135:ILE:HD11	1.91	0.51
1:A:337:SER:HG	3:A:6001:GGA:H372	1.76	0.51
1:B:15:LEU:HD21	1:B:18:TYR:OH	2.10	0.51
1:A:564:LYS:HG2	1:A:565:ASN:N	2.26	0.51
1:A:538:ARG:HG3	1:A:538:ARG:HH11	1.77	0.50
1:A:113:VAL:CG2	1:A:114:ALA:N	2.74	0.50
1:B:527:LYS:HE2	1:B:528:THR:CG2	2.41	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:43:VAL:N	5:B:4183:HOH:O	2.43	0.50
1:B:112:PRO:HB2	1:B:135:ILE:HD13	1.93	0.50
1:A:614:LEU:O	1:A:615:ILE:HD12	2.12	0.50
1:A:228:ASP:OD2	1:A:230:LYS:N	2.45	0.50
1:B:240:ALA:CB	1:B:390:ILE:HG13	2.41	0.50
1:B:137:GLN:HG2	1:B:144:ARG:HB2	1.93	0.49
1:B:131:HIS:HE1	1:B:147:GLU:OE1	1.95	0.49
1:B:137:GLN:CG	1:B:144:ARG:HB2	2.42	0.49
1:B:48:ILE:HD13	1:B:69:LEU:HD21	1.93	0.49
1:B:313:ILE:HD13	1:B:487:PHE:HD2	1.77	0.49
1:B:236:ALA:HB2	1:B:397:GLN:HE22	1.78	0.49
5:A:4487:HOH:O	1:B:160:GLN:HG2	2.12	0.49
1:A:443:ALA:HB3	3:A:6001:GGA:O4	2.12	0.49
1:A:233:LEU:HD21	1:A:398:PHE:HD1	1.77	0.49
1:B:341:GLU:HB3	1:B:611:ILE:HG13	1.94	0.49
1:B:468:HIS:HE1	1:B:470:GLN:HB2	1.76	0.49
1:B:327:LYS:HG3	1:B:515:LEU:HD21	1.93	0.49
1:A:246:ASN:HB2	5:A:4392:HOH:O	2.12	0.49
1:B:601:GLU:H	1:B:601:GLU:CD	2.16	0.49
1:A:459:PHE:HD1	1:A:459:PHE:O	1.96	0.49
1:A:608:GLU:H	1:A:608:GLU:CD	2.15	0.49
1:A:608:GLU:O	1:A:609:SER:C	2.51	0.49
1:A:364:GLY:C	5:A:4456:HOH:O	2.51	0.49
1:B:104:PHE:O	1:B:187:GLY:HA3	2.11	0.49
1:B:233:LEU:HD13	1:B:344:LEU:HD22	1.95	0.49
1:B:215:ARG:HH11	1:B:215:ARG:HG3	1.77	0.49
1:A:82:TRP:CD1	1:A:129:THR:O	2.63	0.48
1:B:120:TRP:NE1	1:B:185:ILE:HD11	2.28	0.48
1:A:311:HIS:CG	5:A:4195:HOH:O	2.58	0.48
1:A:598:ARG:HD2	1:A:611:ILE:HD13	1.94	0.48
1:A:436:ASN:HD22	1:A:436:ASN:C	2.15	0.48
1:B:533:LYS:NZ	1:B:568:GLN:NE2	2.60	0.48
1:A:137:GLN:NE2	1:A:144:ARG:NH1	2.61	0.48
1:B:43:VAL:HG12	1:B:43:VAL:O	2.12	0.48
1:A:366:ILE:HG23	1:A:367:ASN:N	2.27	0.48
1:A:474:ARG:NH1	1:A:476:ILE:HD11	2.29	0.48
1:B:44:PHE:HB3	1:B:55:HIS:CE1	2.49	0.48
1:A:109:PRO:HB2	1:A:197:TYR:CE1	2.48	0.48
1:B:533:LYS:HZ3	1:B:568:GLN:HE22	1.60	0.48
1:B:418:GLU:HG3	1:B:430:GLY:CA	2.43	0.48
1:A:568:GLN:HG2	4:A:3001:ADP:HN61	1.78	0.48
1:B:524:GLU:O	1:B:527:LYS:HG3	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:247:GLN:HG2	5:A:4392:HOH:O	2.13	0.48
1:A:364:GLY:O	1:A:368:GLU:HB2	2.14	0.48
1:B:127:LYS:HB3	1:B:128:ASP:H	1.46	0.48
1:B:262:GLU:OE1	1:B:262:GLU:HA	2.14	0.47
1:A:165:LEU:CD2	1:A:179:THR:HG23	2.44	0.47
1:A:287:LEU:HA	1:A:290:LEU:HD12	1.96	0.47
1:B:196:GLU:O	1:B:197:TYR:HB2	2.13	0.47
1:B:150:VAL:HG12	1:B:151:ILE:HD12	1.96	0.47
1:B:581:ILE:CG2	1:B:597:LEU:HD22	2.44	0.47
1:B:150:VAL:C	1:B:151:ILE:HD12	2.34	0.47
1:A:538:ARG:HH11	1:A:538:ARG:CG	2.28	0.47
1:B:427:LEU:HD21	1:B:478:VAL:HG22	1.97	0.47
1:B:95:ASN:O	1:B:97:ASN:N	2.48	0.47
1:B:274:TYR:CE1	1:B:504:LEU:HD11	2.49	0.47
1:B:547:SER:OG	1:B:551:GLU:HG2	2.15	0.47
1:A:522:ASN:O	1:A:526:VAL:HG23	2.14	0.47
1:A:575:LYS:HE2	5:A:4254:HOH:O	2.15	0.46
1:A:274:TYR:CZ	1:A:504:LEU:HD11	2.50	0.46
1:B:316:ARG:HD2	1:B:584:CYS:HB3	1.97	0.46
1:B:400:GLU:HG3	1:B:410:THR:OG1	2.15	0.46
1:A:98:ILE:HD13	1:A:276:HIS:CD2	2.50	0.46
1:B:104:PHE:CE2	1:B:110:ARG:HG2	2.49	0.46
1:B:347:GLU:HA	1:B:347:GLU:OE1	2.15	0.46
1:B:85:PHE:HA	1:B:189:MET:CE	2.45	0.46
1:A:137:GLN:HE22	1:A:144:ARG:NH1	2.13	0.46
1:B:339:HIS:N	1:B:339:HIS:CD2	2.84	0.46
1:B:121:ASP:O	1:B:130:GLY:HA2	2.16	0.46
1:A:421:TRP:HE3	1:A:471:ASN:HD22	1.64	0.46
1:B:131:HIS:CG	1:B:132:VAL:N	2.83	0.46
1:B:48:ILE:CG2	1:B:68:PHE:HE2	2.29	0.46
1:B:240:ALA:HB2	1:B:390:ILE:HG13	1.98	0.46
1:A:547:SER:HB3	1:A:553:LEU:HD11	1.97	0.46
1:B:418:GLU:HG3	1:B:430:GLY:HA3	1.98	0.46
1:B:117:LEU:O	1:B:188:TRP:HA	2.16	0.46
1:B:552:VAL:HG11	1:B:555:LYS:HE3	1.98	0.46
1:B:379:ARG:HE	1:B:485:LEU:HB2	1.81	0.46
1:B:221:GLN:OE1	1:B:250:TYR:HE2	1.99	0.46
1:A:138:LEU:HD12	1:A:143:VAL:CG1	2.42	0.46
1:B:85:PHE:HA	1:B:189:MET:SD	2.56	0.46
1:B:96:ASP:CG	1:B:309:ARG:HH21	2.18	0.46
1:B:66:PHE:CE2	1:B:134:ILE:HG21	2.51	0.46
1:A:268:ASN:ND2	1:A:592:TYR:H	2.11	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:324:ARG:NH2	1:B:572:CYS:O	2.45	0.45
1:B:348:ARG:NH1	1:B:351:GLU:OE2	2.48	0.45
1:A:313:ILE:HD13	1:A:487:PHE:HD2	1.81	0.45
1:B:459:PHE:CD2	1:B:459:PHE:N	2.82	0.45
1:A:548:HIS:C	1:A:550:GLU:H	2.19	0.45
1:A:174:TYR:CD1	1:A:174:TYR:N	2.84	0.45
1:A:498:LYS:HB2	1:A:535:ILE:HA	1.99	0.45
1:B:303:LEU:O	1:B:307:ARG:HG2	2.17	0.45
1:B:379:ARG:H	1:B:379:ARG:HD3	1.82	0.45
1:B:449:ASP:OD1	1:B:452:ARG:NH2	2.40	0.45
1:B:608:GLU:O	1:B:609:SER:C	2.54	0.45
1:A:379:ARG:NH1	1:A:436:ASN:HB2	2.29	0.45
1:B:386:GLN:NE2	1:B:387:ASP:O	2.50	0.44
3:A:6001:GGA:H491	5:A:4176:HOH:O	2.17	0.44
1:B:120:TRP:NE1	1:B:185:ILE:CD1	2.80	0.44
1:B:138:LEU:HD12	1:B:143:VAL:CG1	2.43	0.44
1:B:112:PRO:HB2	1:B:135:ILE:CD1	2.48	0.44
1:A:204:ILE:HG21	1:A:262:GLU:HG2	1.99	0.44
1:A:56:LYS:HA	1:A:57:TRP:HA	1.61	0.44
1:B:444:TRP:HB2	1:B:493:VAL:HG12	1.99	0.44
1:B:28:SER:HB2	1:B:57:TRP:O	2.17	0.44
1:B:387:ASP:HB2	3:B:6002:GGA:HN81	1.81	0.44
1:A:25:ILE:HG12	1:A:62:PHE:CE1	2.52	0.44
1:A:533:LYS:NZ	1:A:568:GLN:NE2	2.66	0.44
1:A:300:ARG:NH1	5:A:4196:HOH:O	2.50	0.44
1:B:468:HIS:HE1	1:B:470:GLN:HE21	1.66	0.44
1:A:346:LEU:HD23	1:A:346:LEU:HA	1.80	0.44
1:A:63:ALA:O	1:A:67:LEU:HG	2.18	0.44
1:B:584:CYS:HB2	1:B:596:CYS:SG	2.58	0.44
1:A:430:GLY:O	1:A:432:GLY:N	2.49	0.44
1:B:360:ASN:OD1	1:B:361:PRO:HD2	2.18	0.44
1:A:10:ALA:N	5:A:4348:HOH:O	2.50	0.44
1:A:151:ILE:HD12	1:A:162:THR:CB	2.44	0.44
1:B:56:LYS:HA	1:B:57:TRP:HA	1.66	0.44
1:B:311:HIS:O	1:B:313:ILE:HD12	2.18	0.44
1:A:137:GLN:CG	1:A:144:ARG:HB2	2.46	0.43
1:A:421:TRP:HA	1:A:426:GLN:O	2.18	0.43
1:A:421:TRP:HE3	1:A:471:ASN:ND2	2.16	0.43
1:A:430:GLY:C	1:A:432:GLY:H	2.21	0.43
1:A:68:PHE:HE1	1:A:93:VAL:HG11	1.82	0.43
1:B:521:VAL:HG22	1:B:567:TYR:CD1	2.53	0.43
1:B:535:ILE:HG23	1:B:564:LYS:O	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:109:PRO:HA	1:A:172:GLY:O	2.17	0.43
1:A:411:ARG:HA	5:A:4313:HOH:O	2.19	0.43
1:A:109:PRO:HB2	1:A:197:TYR:CD1	2.53	0.43
1:B:120:TRP:HB2	1:B:131:HIS:HB3	1.99	0.43
1:B:20:PRO:HD3	1:B:160:GLN:O	2.17	0.43
1:A:308:ARG:HH11	1:A:308:ARG:HG3	1.83	0.43
1:B:512:ARG:HH11	1:B:512:ARG:HG2	1.82	0.43
1:B:444:TRP:CB	1:B:493:VAL:HG12	2.49	0.43
1:A:74:VAL:HG22	1:A:75:PHE:N	2.34	0.43
1:B:48:ILE:O	1:B:49:ASP:CB	2.66	0.43
1:A:205:ALA:O	1:A:207:GLU:N	2.52	0.43
1:B:337:SER:O	1:B:338:CYS:CB	2.58	0.43
1:A:201:GLN:CG	1:A:202:PRO:HD2	2.47	0.43
1:A:15:LEU:HD21	1:A:18:TYR:CE1	2.53	0.43
1:A:348:ARG:HD2	1:A:351:GLU:OE1	2.19	0.43
1:B:131:HIS:CE1	1:B:150:VAL:HG21	2.54	0.43
1:A:10:ALA:N	1:A:152:HIS:O	2.52	0.43
1:B:490:LEU:O	1:B:493:VAL:HG13	2.18	0.43
1:B:314:THR:HA	1:B:587:THR:O	2.18	0.43
1:A:234:GLN:O	1:A:237:TYR:HB3	2.19	0.42
1:A:481:ARG:NH2	5:A:4229:HOH:O	2.52	0.42
1:B:45:ARG:O	1:B:56:LYS:HE3	2.19	0.42
1:A:468:HIS:HA	1:A:469:PRO:HD2	1.85	0.42
1:B:503:ILE:O	1:B:506:SER:HB3	2.19	0.42
1:A:62:PHE:HB2	1:A:148:GLN:HB2	2.01	0.42
1:B:522:ASN:ND2	1:B:522:ASN:C	2.68	0.42
1:B:44:PHE:O	1:B:56:LYS:HG3	2.20	0.42
1:A:107:GLY:CA	1:A:174:TYR:O	2.66	0.42
1:A:337:SER:O	1:A:338:CYS:CB	2.60	0.42
1:A:245:ILE:HG23	1:A:345:ILE:HG21	2.01	0.42
1:B:533:LYS:HZ3	1:B:568:GLN:NE2	2.16	0.42
1:A:88:ARG:HG3	1:A:201:GLN:HB2	2.02	0.42
1:A:317:MET:O	1:A:584:CYS:HA	2.19	0.42
1:A:141:ASN:HD22	1:A:141:ASN:H	1.67	0.42
1:B:431:GLU:HA	1:B:431:GLU:OE2	2.19	0.42
1:A:536:ALA:H	1:A:563:GLN:NE2	2.18	0.42
1:A:168:VAL:CG2	1:A:175:THR:HB	2.49	0.42
1:B:111:ALA:HB2	1:B:174:TYR:CZ	2.55	0.42
1:A:334:ASP:OD2	1:A:444:TRP:HB2	2.20	0.42
1:B:532:VAL:HG22	1:B:567:TYR:CD1	2.55	0.42
1:A:582:GLN:NE2	4:A:3001:ADP:O3'	2.53	0.42
1:A:501:LEU:HD23	1:A:504:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:339:HIS:HB3	1:A:398:PHE:CE2	2.55	0.42
1:B:601:GLU:N	1:B:601:GLU:CD	2.73	0.42
1:B:219:LYS:HE3	1:B:219:LYS:HB3	1.89	0.42
1:A:465:ARG:HD3	1:A:474:ARG:NH2	2.34	0.42
1:A:108:SER:HA	1:A:109:PRO:HD3	1.93	0.42
1:A:137:GLN:CG	1:A:144:ARG:HD2	2.50	0.41
1:B:151:ILE:N	1:B:151:ILE:HD12	2.35	0.41
1:A:16:LEU:HD21	1:A:27:SER:HB2	2.02	0.41
1:B:244:VAL:CG1	1:B:245:ILE:N	2.83	0.41
1:A:382:VAL:HB	1:A:410:THR:HG22	2.00	0.41
1:B:276:HIS:CE1	1:B:280:LYS:HE2	2.55	0.41
1:A:324:ARG:NH2	1:A:572:CYS:O	2.49	0.41
1:A:576:VAL:HG12	1:A:577:ASP:OD1	2.20	0.41
1:B:381:PHE:CD1	1:B:409:GLU:HB2	2.56	0.41
1:B:418:GLU:CG	1:B:418:GLU:O	2.68	0.41
1:B:245:ILE:HG13	1:B:246:ASN:OD1	2.20	0.41
1:A:379:ARG:HH11	1:A:379:ARG:HG3	1.85	0.41
1:B:515:LEU:HD12	1:B:568:GLN:OE1	2.21	0.41
1:B:131:HIS:CG	1:B:132:VAL:H	2.39	0.41
1:A:440:LYS:HE2	1:A:444:TRP:CD1	2.55	0.41
1:B:29:ASP:O	1:B:30:TYR:CG	2.73	0.41
1:B:231:ASP:CB	1:B:234:GLN:HE21	2.33	0.41
1:A:268:ASN:ND2	1:A:591:ASN:HA	2.36	0.41
1:B:119:ILE:HD12	1:B:119:ILE:N	2.36	0.41
1:A:237:TYR:CE2	1:A:345:ILE:HG12	2.55	0.41
1:A:18:TYR:CG	1:B:466:THR:HG21	2.56	0.41
1:A:598:ARG:HA	1:A:610:ASP:O	2.21	0.41
1:A:498:LYS:C	1:A:500:ILE:H	2.24	0.41
1:B:204:ILE:HD11	1:B:208:LEU:HD12	2.03	0.41
1:A:253:TYR:CZ	1:A:581:ILE:HD11	2.56	0.41
1:A:313:ILE:HD13	1:A:487:PHE:CD2	2.56	0.41
1:B:470:GLN:HB2	1:B:470:GLN:HE21	1.71	0.41
1:B:196:GLU:HG2	1:B:197:TYR:CD1	2.55	0.41
1:A:405:GLN:C	1:A:407:GLY:H	2.24	0.41
1:B:312:MET:HA	1:B:488:GLU:OE2	2.20	0.41
1:A:527:LYS:O	1:A:548:HIS:CB	2.66	0.40
1:B:144:ARG:HD3	1:B:161:TRP:CG	2.56	0.40
1:A:216:LEU:HD13	1:A:353:GLY:O	2.21	0.40
1:A:15:LEU:HD21	1:A:18:TYR:OH	2.20	0.40
1:B:564:LYS:HG2	1:B:565:ASN:N	2.36	0.40
1:A:462:VAL:HG11	1:A:480:LEU:HD12	2.03	0.40
1:B:560:PHE:N	1:B:560:PHE:CD1	2.89	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:531:ALA:HA	1:A:544:ASP:O	2.21	0.40
1:B:597:LEU:HD11	1:B:614:LEU:HD13	2.03	0.40
1:B:322:ASP:OD2	1:B:322:ASP:C	2.60	0.40
1:A:156:PRO:O	1:A:159:GLN:HB2	2.21	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	581/619 (94%)	534 (92%)	40 (7%)	7 (1%)	19	54
1	B	581/619 (94%)	526 (90%)	47 (8%)	8 (1%)	16	49
All	All	1162/1238 (94%)	1060 (91%)	87 (8%)	15 (1%)	18	51

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	THR
1	A	127	LYS
1	A	334	ASP
1	B	96	ASP
1	B	127	LYS
1	B	334	ASP
1	A	431	GLU
1	B	49	ASP
1	B	338	CYS
1	A	49	ASP
1	A	338	CYS
1	B	95	ASN
1	B	365	LEU
1	B	522	ASN
1	A	206	GLY

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	505/530 (95%)	489 (97%)	16 (3%)	51	85
1	B	505/530 (95%)	481 (95%)	24 (5%)	35	72
All	All	1010/1060 (95%)	970 (96%)	40 (4%)	42	79

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	128	ASP
1	A	141	ASN
1	A	149	ASN
1	A	184	THR
1	A	245	ILE
1	A	300	ARG
1	A	355	LYS
1	A	390	ILE
1	A	436	ASN
1	A	459	PHE
1	A	470	GLN
1	A	471	ASN
1	A	481	ARG
1	A	523	ASP
1	A	608	GLU
1	B	91	ARG
1	B	127	LYS
1	B	128	ASP
1	B	201	GLN
1	B	204	ILE
1	B	233	LEU
1	B	247	GLN
1	B	261	GLN
1	B	352	GLN
1	B	377	ARG
1	B	379	ARG
1	B	386	GLN
1	B	389	ASP

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Mol	Chain	Res	Type
1	B	421	TRP
1	B	426	GLN
1	B	459	PHE
1	B	466	THR
1	B	470	GLN
1	B	471	ASN
1	B	522	ASN
1	B	524	GLU
1	B	527	LYS
1	B	560	PHE
1	B	615	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	70	ASN
1	A	95	ASN
1	A	131	HIS
1	A	137	GLN
1	A	139	HIS
1	A	141	ASN
1	A	149	ASN
1	A	152	HIS
1	A	268	ASN
1	A	310	HIS
1	A	339	HIS
1	A	352	GLN
1	A	395	HIS
1	A	405	GLN
1	A	436	ASN
1	A	470	GLN
1	A	471	ASN
1	A	563	GLN
1	A	568	GLN
1	B	55	HIS
1	B	102	GLN
1	B	131	HIS
1	B	157	GLN
1	B	201	GLN
1	B	234	GLN
1	B	247	GLN

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Mol	Chain	Res	Type
1	B	251	HIS
1	B	268	ASN
1	B	286	ASN
1	B	310	HIS
1	B	339	HIS
1	B	357	ASN
1	B	367	ASN
1	B	375	HIS
1	B	395	HIS
1	B	397	GLN
1	B	468	HIS
1	B	470	GLN
1	B	471	ASN
1	B	522	ASN
1	B	563	GLN
1	B	568	GLN
1	B	582	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	ADP	A	3001	2	29,29,29	1.69	2 (6%)	45,45,45	1.47	5 (11%)
3	GGA	A	6001	2	32,32,32	5.35	9 (28%)	40,42,42	3.02	13 (32%)
4	ADP	B	3002	2	29,29,29	1.77	2 (6%)	45,45,45	1.51	7 (15%)
3	GGA	B	6002	2	32,32,32	5.80	8 (25%)	40,42,42	3.55	14 (35%)

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	ADP	A	3001	2	-	0/16/32/32	0/1/3/3
3	GGA	A	6001	2	-	0/37/40/40	0/0/0/0
4	ADP	B	3002	2	-	0/16/32/32	0/1/3/3
3	GGA	B	6002	2	-	0/37/40/40	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	6002	GGA	P34-C29	-23.97	1.62	1.82
3	A	6001	GGA	P34-C29	-20.63	1.65	1.82
3	B	6002	GGA	P34-C33	-17.98	1.64	1.79
3	A	6001	GGA	P34-C33	-17.98	1.64	1.79
3	B	6002	GGA	P34-O56	-8.68	1.32	1.49
3	A	6001	GGA	P34-O56	-8.21	1.33	1.49
4	B	3002	ADP	O3'-C3'	-7.71	1.24	1.43
4	A	3001	ADP	O3'-C3'	-7.53	1.24	1.43
3	B	6002	GGA	O2-C1	6.05	1.42	1.22
3	B	6002	GGA	P60-O64	-5.42	1.34	1.54
3	A	6001	GGA	P60-O64	-5.34	1.35	1.54
3	A	6001	GGA	O2-C1	4.54	1.37	1.22
3	A	6001	GGA	O4-C1	4.02	1.44	1.30
3	A	6001	GGA	C29-N30	-3.49	1.41	1.45
4	B	3002	ADP	C2-N1	3.08	1.40	1.33
3	B	6002	GGA	C14-N18	2.99	1.40	1.34
4	A	3001	ADP	C2-N1	2.99	1.39	1.33
3	B	6002	GGA	O4-C1	-2.86	1.20	1.30
3	B	6002	GGA	C29-N30	-2.86	1.42	1.45
3	A	6001	GGA	C14-N18	2.40	1.39	1.34
3	A	6001	GGA	P60-O58	-2.03	1.56	1.60

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	6002	GGA	O58-P34-C29	15.61	118.07	107.93
3	A	6001	GGA	O58-P34-C29	13.29	116.56	107.93
3	B	6002	GGA	P34-C29-N30	-9.49	101.62	112.26
3	A	6001	GGA	P34-C29-N30	-8.17	103.09	112.26
3	B	6002	GGA	O4-C1-C6	6.41	131.26	116.88
4	B	3002	ADP	N3-C2-N1	-6.30	123.44	128.71
4	A	3001	ADP	N3-C2-N1	-5.98	123.71	128.71
3	B	6002	GGA	C20-N18-C14	4.33	127.69	121.52
3	A	6001	GGA	O56-P34-C29	4.16	110.81	108.54
3	B	6002	GGA	C22-C20-N18	4.08	118.15	110.31
3	A	6001	GGA	C22-C20-N18	4.05	118.08	110.31
3	A	6001	GGA	O56-P34-C33	-3.52	100.22	110.75
3	A	6001	GGA	C20-N18-C14	3.49	126.50	121.52
3	B	6002	GGA	O56-P34-C33	-3.43	100.48	110.75
4	A	3001	ADP	C4'-O4'-C1'	3.24	113.27	109.75
3	B	6002	GGA	O26-C24-C20	-3.22	112.81	120.42
4	B	3002	ADP	C4'-O4'-C1'	3.10	113.12	109.75
3	B	6002	GGA	C12-C10-C6	-3.01	108.81	114.43
3	A	6001	GGA	O26-C24-C20	-2.86	113.67	120.42
3	B	6002	GGA	O2-C1-C6	-2.84	110.17	118.36
3	B	6002	GGA	C20-C24-N30	2.80	122.75	116.52
4	A	3001	ADP	C3'-C2'-C1'	2.75	105.20	100.91
3	A	6001	GGA	C37-C39-N44	-2.60	105.18	112.09
3	A	6001	GGA	O2-C1-C6	2.56	125.75	118.36
3	B	6002	GGA	O4-C1-O2	-2.51	118.39	124.07
3	B	6002	GGA	C37-C39-N44	-2.43	105.63	112.09
4	A	3001	ADP	C2-N1-C6	2.42	123.14	118.77
4	B	3002	ADP	C2-N1-C6	2.39	123.09	118.77
3	B	6002	GGA	C29-P34-C33	2.31	111.30	106.86
4	A	3001	ADP	C5-C6-N6	2.27	125.85	120.72
3	A	6001	GGA	C12-C10-C6	-2.24	110.25	114.43
3	A	6001	GGA	O4-C1-C6	-2.23	111.87	116.88
3	A	6001	GGA	C29-P34-C33	2.23	111.14	106.86
3	B	6002	GGA	O56-P34-C29	2.22	109.75	108.54
4	B	3002	ADP	C3'-C2'-C1'	2.14	104.26	100.91
4	B	3002	ADP	C8-N9-C4	-2.12	105.28	106.90
4	B	3002	ADP	C4-C5-N7	2.12	111.33	109.52
3	A	6001	GGA	C20-C24-N30	2.06	121.11	116.52
4	B	3002	ADP	C5-C6-N6	2.02	125.30	120.72

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

## 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	589/619 (95%)	-0.42	6 (1%) 79 79	7, 27, 48, 63	0
1	B	589/619 (95%)	-0.37	5 (0%) 83 83	8, 29, 48, 67	0
All	All	1178/1238 (95%)	-0.40	11 (0%) 81 81	7, 28, 48, 67	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	43	VAL	3.5
1	A	44	PHE	3.1
1	A	172	GLY	3.1
1	B	562	GLU	2.9
1	B	31	SER	2.8
1	A	453	GLU	2.5
1	B	29	ASP	2.3
1	A	157	GLN	2.1
1	A	29	ASP	2.1
1	B	141	ASN	2.1
1	B	171	ASN	2.0

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

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GGA	B	6002	33/33	0.20	1.27	22,37,54,56	0
3	GGA	A	6001	33/33	0.21	1.03	22,31,45,49	0
4	ADP	B	3002	27/27	0.14	-0.22	6,18,20,21	0
2	MG	A	5001	1/1	0.15	-0.45	8,8,8,8	0
4	ADP	A	3001	27/27	0.12	-0.47	1,18,21,23	0
2	MG	B	5003	1/1	0.12	-0.73	10,10,10,10	0
2	MG	A	5002	1/1	0.15	-1.30	8,8,8,8	0
2	MG	B	5004	1/1	0.12	-3.21	12,12,12,12	0

## 6.5 Other polymers

There are no such residues in this entry.