



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 06:02 am BST

PDB ID : 1ASL  
Title : CRYSTAL STRUCTURES OF ESCHERICHIA COLI ASPARTATE AMINO-TRANSFERASE IN TWO CONFORMATIONS: COMPARISON OF AN UNLIGANDED OPEN AND TWO LIGANDED CLOSED FORMS  
Authors : Jaeger, J.; Jansonius, J.N.  
Deposited on : 1993-09-16  
Resolution : 2.60 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

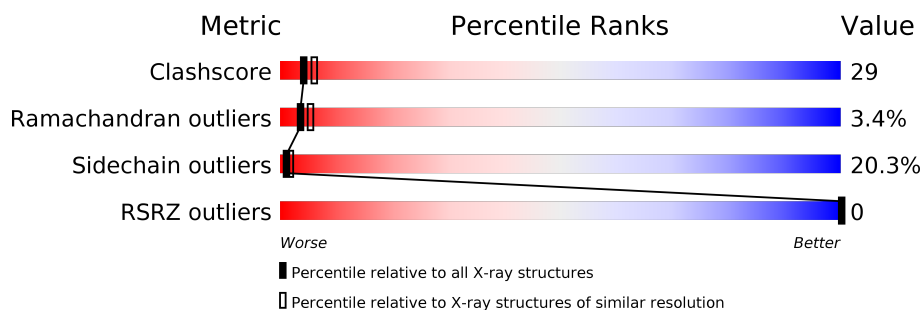
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

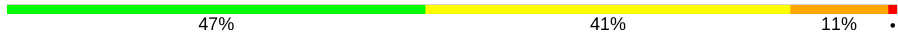
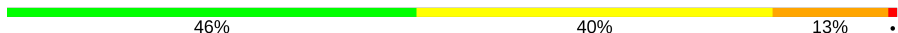
The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	396	 47% 41% 11% .
1	B	396	 46% 40% 13% .

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
2	PLA	B	410	-	-	X	-

## 2 Entry composition [i](#)

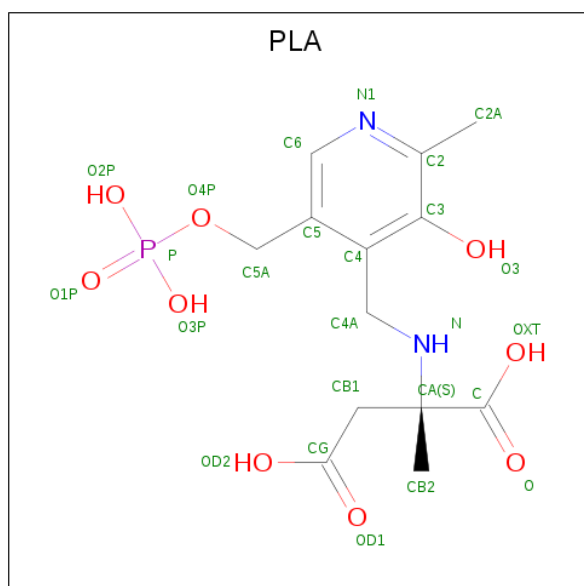
There are 3 unique types of molecules in this entry. The entry contains 6407 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			3069	1936	536	584	13			
1	B	396	Total	C	N	O	S	0	0	0
			3069	1936	536	584	13			

- Molecule 2 is 2-[(3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YLMETHYL)-AMINO]-2-METHYL-SUCCINIC ACID (three-letter code: PLA) (formula:  $C_{13}H_{19}N_2O_9P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	13	2	9	1		
2	B	1	Total	C	N	O	P	0	0
			25	13	2	9	1		

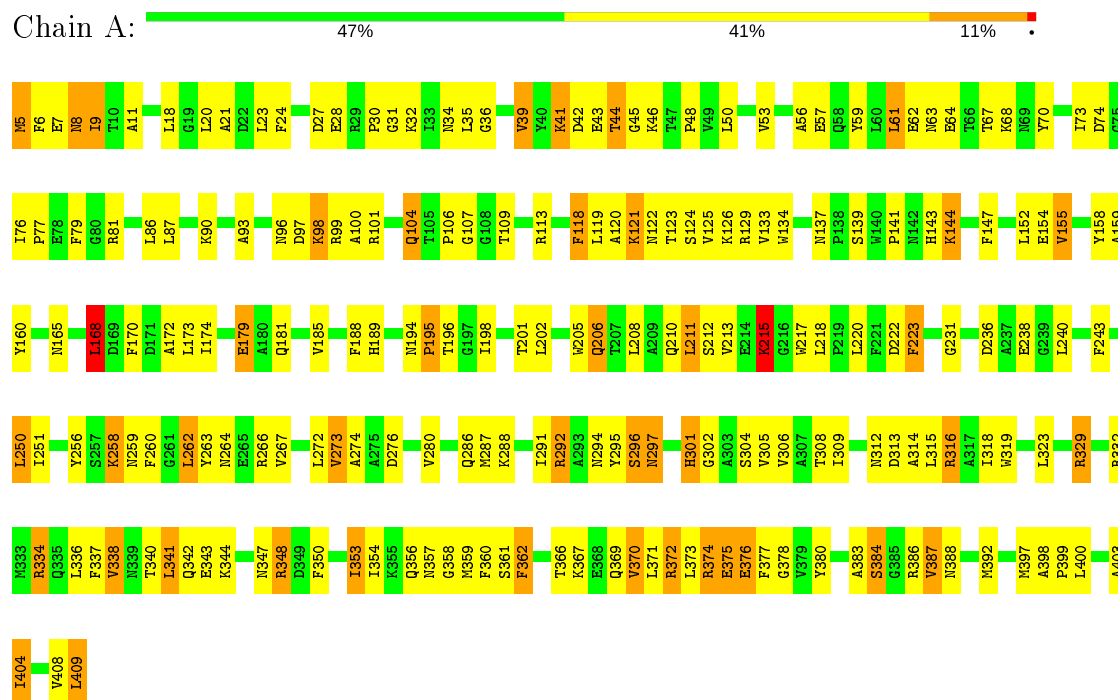
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	113	Total 113	O 113	0	0
3	B	106	Total 106	O 106	0	0

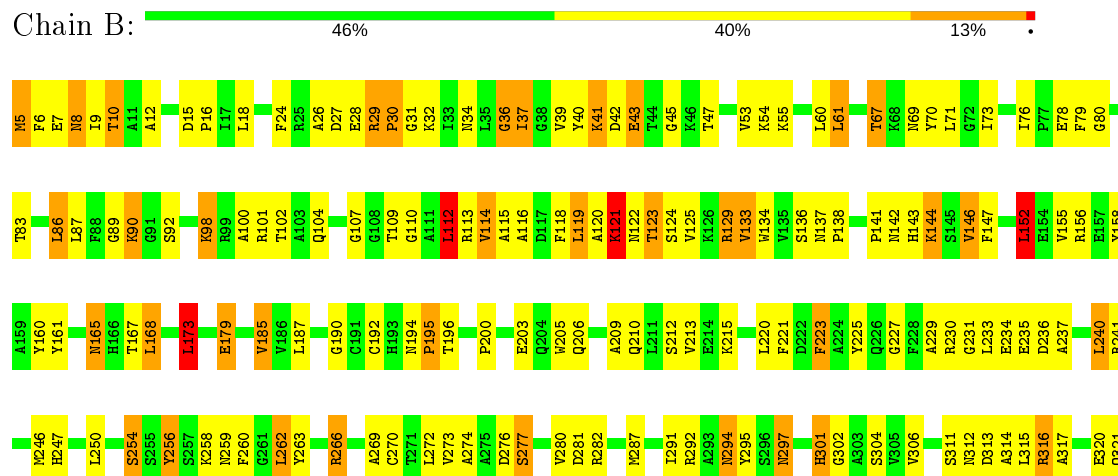
### 3 Residue-property plots

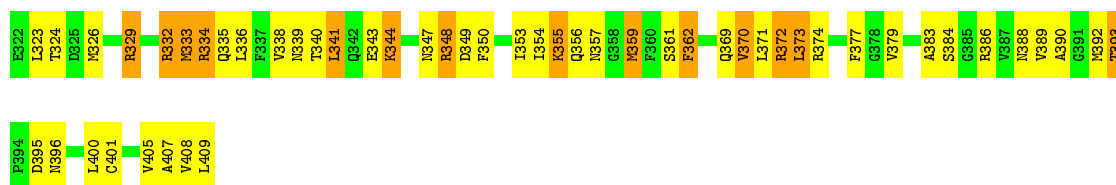
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

#### • Molecule 1: ASPARTATE AMINOTRANSFERASE



#### • Molecule 1: ASPARTATE AMINOTRANSFERASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.60 Å 78.80 Å 89.60 Å 90.00° 118.60° 90.00°	Depositor
Resolution (Å)	10.00 – 2.60 10.98 – 2.51	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.60) 91.2 (10.98-2.51)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 2.51 Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.183 , (Not available) 0.181 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.0	Xtriage
Anisotropy	0.524	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.23 , 78.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h-l 0.000 for -h-l,k,h 0.457 for h,-k,-h-l 0.000 for l,-k,h 0.000 for -h-l,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6407	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.84	0/3130	1.06	5/4240 (0.1%)
1	B	0.83	0/3130	1.07	7/4240 (0.2%)
All	All	0.83	0/6260	1.06	12/8480 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	409	LEU	CA-CB-CG	9.29	136.66	115.30
1	A	168	LEU	CA-CB-CG	8.81	135.57	115.30
1	B	89	GLY	N-CA-C	7.60	132.11	113.10
1	B	250	LEU	CA-CB-CG	6.67	130.64	115.30
1	A	334	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	B	359	MET	CG-SD-CE	-6.46	89.86	100.20
1	B	173	LEU	CA-CB-CG	5.86	128.78	115.30
1	A	74	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	316	ARG	NE-CZ-NH2	-5.37	117.62	120.30
1	B	27	ASP	N-CA-C	-5.30	96.70	111.00
1	B	377	PHE	N-CA-C	5.25	125.17	111.00
1	B	112	LEU	CA-CB-CG	5.15	127.15	115.30

There are no chirality outliers.



All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	256	TYR	Sidechain

## 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	3069	0	3018	174	0
1	B	3069	0	3018	209	0
2	A	25	0	14	5	0
2	B	25	0	14	10	0
3	A	113	0	0	9	0
3	B	106	0	0	15	0
All	All	6407	0	6064	361	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 29.

All (361) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:PHE:HB3	1:B:262:LEU:HD22	1.34	1.07
1:A:39:VAL:HG21	1:B:69:ASN:HD22	1.21	1.03
1:A:260:PHE:HB3	1:A:262:LEU:HD22	1.44	1.00
1:A:196:THR:HA	1:A:362:PHE:HB2	1.43	1.00
1:B:8:ASN:H	1:B:8:ASN:HD22	1.12	0.98
1:A:39:VAL:HG21	1:B:69:ASN:ND2	1.82	0.94
1:A:294:ASN:HD21	1:B:294:ASN:HD21	0.95	0.94
1:A:61:LEU:HD12	1:B:61:LEU:HD12	1.51	0.90
1:A:210:GLN:O	1:A:213:VAL:HG12	1.73	0.89
1:B:258:LYS:NZ	2:B:410:PLA:H4A2	1.88	0.88
1:A:196:THR:HG22	1:A:362:PHE:CD2	2.09	0.87
1:B:123:THR:CG2	1:B:125:VAL:HG13	2.07	0.85
1:A:8:ASN:H	1:A:8:ASN:HD22	1.23	0.84
1:A:296:SER:HB2	1:B:109:THR:HG21	1.59	0.83
1:B:107:GLY:HA3	1:B:266:ARG:NH1	1.92	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:LEU:HD12	1:B:326:MET:HE3	1.61	0.82
1:A:294:ASN:ND2	1:B:294:ASN:HD21	1.77	0.81
1:A:6:PHE:HD1	1:A:9:ILE:HD11	1.46	0.81
1:A:276:ASP:O	1:A:280:VAL:HG23	1.81	0.80
1:B:8:ASN:H	1:B:8:ASN:ND2	1.75	0.80
1:B:258:LYS:HZ1	2:B:410:PLA:H4A2	1.44	0.80
1:A:294:ASN:HD21	1:B:294:ASN:ND2	1.79	0.80
1:B:129:ARG:HD2	1:B:134:TRP:HE1	1.47	0.80
1:B:121:LYS:HD3	1:B:121:LYS:O	1.83	0.79
1:A:350:PHE:HB3	1:A:353:ILE:HD13	1.66	0.78
1:B:338:VAL:HG11	1:B:354:ILE:CD1	2.14	0.78
1:B:209:ALA:O	1:B:213:VAL:HG23	1.83	0.78
1:A:120:ALA:HB2	1:A:152:LEU:HD13	1.65	0.78
1:A:6:PHE:O	1:A:9:ILE:HG13	1.84	0.78
1:A:107:GLY:HA3	1:A:266:ARG:NH1	2.01	0.76
1:A:366:THR:HG22	1:A:369:GLN:H	1.48	0.76
1:A:122:ASN:OD1	1:A:286:GLN:HG2	1.85	0.76
1:A:274:ALA:HB3	1:A:280:VAL:HG22	1.66	0.76
1:B:40:TYR:CE2	1:B:326:MET:HG2	2.20	0.75
1:B:18:LEU:HD21	3:B:515:HOH:O	1.86	0.74
2:B:410:PLA:H4A1	3:B:509:HOH:O	1.87	0.74
1:A:129:ARG:NH2	1:A:181:GLN:HG3	2.03	0.73
1:B:110:GLY:O	1:B:114:VAL:HG13	1.87	0.73
1:B:355:LYS:HD3	1:B:355:LYS:O	1.89	0.73
1:B:227:GLY:HA2	1:B:231:GLY:O	1.88	0.72
1:A:374:ARG:HG3	1:A:375:GLU:N	2.05	0.72
1:B:338:VAL:HG11	1:B:354:ILE:HD11	1.72	0.72
1:A:77:PRO:O	1:A:81:ARG:HG3	1.90	0.71
1:A:212:SER:HB3	1:A:217:TRP:HE3	1.56	0.70
1:A:120:ALA:HB2	1:A:152:LEU:CD1	2.21	0.70
1:B:223:PHE:CE2	1:B:240:LEU:HD12	2.27	0.70
1:A:99:ARG:O	1:A:280:VAL:HG21	1.91	0.70
1:B:274:ALA:HB3	1:B:280:VAL:HB	1.74	0.69
1:A:286:GLN:OE1	1:B:12:ALA:HB2	1.92	0.69
1:B:165:ASN:HD21	1:B:167:THR:HG22	1.57	0.69
1:A:123:THR:OG1	1:A:125:VAL:HG23	1.92	0.68
1:A:11:ALA:HB2	1:B:282:ARG:HG2	1.75	0.68
1:A:159:ALA:O	1:A:173:LEU:HD23	1.94	0.68
1:B:8:ASN:N	1:B:8:ASN:HD22	1.90	0.68
1:B:350:PHE:HB3	1:B:353:ILE:HD13	1.74	0.68
1:A:266:ARG:HD2	1:B:297:ASN:O	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:LYS:HG2	1:B:359:MET:HE3	1.76	0.67
1:B:76:ILE:O	1:B:79:PHE:HB3	1.94	0.67
1:B:195:PRO:HB3	1:B:386:ARG:HG3	1.77	0.67
1:B:129:ARG:NH1	1:B:129:ARG:HB3	2.09	0.67
1:B:142:ASN:O	1:B:146:VAL:HG13	1.95	0.67
1:B:165:ASN:ND2	1:B:167:THR:HG22	2.10	0.66
1:B:67:THR:HG21	3:B:516:HOH:O	1.94	0.65
1:A:329:ARG:NH2	1:A:392:MET:O	2.29	0.65
1:A:168:LEU:O	1:A:168:LEU:HD13	1.95	0.65
1:A:297:ASN:O	1:B:266:ARG:HD2	1.97	0.65
1:B:98:LYS:O	1:B:101:ARG:NH1	2.30	0.64
1:B:133:VAL:HB	1:B:185:VAL:HG22	1.78	0.64
1:B:329:ARG:NH2	1:B:392:MET:O	2.30	0.63
1:A:266:ARG:NH2	2:A:410:PLA:O1P	2.32	0.63
1:A:73:ILE:CG2	1:A:288:LYS:HG2	2.28	0.63
1:B:260:PHE:HB3	1:B:262:LEU:CD2	2.21	0.63
1:A:24:PHE:CE1	1:A:34:ASN:HB2	2.33	0.63
1:B:40:TYR:OH	1:B:326:MET:HA	1.99	0.63
1:A:109:THR:O	1:A:109:THR:HG22	1.98	0.63
1:A:220:LEU:HD12	1:A:251:ILE:HB	1.80	0.63
1:B:374:ARG:HD3	3:B:490:HOH:O	1.99	0.63
1:A:398:ALA:H	1:A:399:PRO:HD2	1.64	0.62
1:A:134:TRP:HH2	1:A:179:GLU:HG2	1.64	0.62
1:A:46:LYS:O	1:A:48:PRO:HD3	1.99	0.62
1:A:286:GLN:NE2	1:B:10:THR:O	2.32	0.62
1:B:6:PHE:O	1:B:9:ILE:HG13	1.99	0.62
1:A:238:GLU:HB2	3:A:413:HOH:O	2.00	0.62
1:B:266:ARG:HH22	2:B:410:PLA:P	2.21	0.62
1:A:42:ASP:O	1:A:45:GLY:N	2.31	0.62
1:A:212:SER:HB3	1:A:217:TRP:CE3	2.35	0.62
1:A:76:ILE:O	1:A:79:PHE:HB3	1.99	0.62
1:B:338:VAL:HG11	1:B:354:ILE:HD12	1.81	0.61
1:B:333:MET:CE	1:B:336:LEU:HD23	2.30	0.61
1:B:137:ASN:O	1:B:160:TYR:HB3	2.01	0.61
1:A:179:GLU:HG3	1:A:179:GLU:O	1.99	0.61
1:A:101:ARG:HD3	3:A:491:HOH:O	2.00	0.61
1:A:70:TYR:HE1	1:B:266:ARG:HH21	1.48	0.61
1:A:8:ASN:ND2	1:A:8:ASN:H	1.97	0.60
1:B:323:LEU:HA	1:B:326:MET:CE	2.31	0.60
2:B:410:PLA:N	2:B:410:PLA:O3	2.29	0.60
1:B:24:PHE:CZ	1:B:32:LYS:HG3	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:348:ARG:NH1	1:B:409:LEU:HB3	2.17	0.59
1:A:334:ARG:NH2	3:A:513:HOH:O	2.34	0.59
1:B:196:THR:HA	1:B:362:PHE:HB2	1.85	0.59
1:B:123:THR:HG22	1:B:125:VAL:H	1.67	0.59
1:A:73:ILE:HD11	1:B:18:LEU:HD12	1.84	0.59
1:B:254:SER:HB2	3:B:492:HOH:O	2.02	0.59
1:B:258:LYS:HG2	1:B:359:MET:CE	2.33	0.59
1:A:348:ARG:HB3	1:A:350:PHE:CE1	2.38	0.58
1:A:356:GLN:C	1:A:357:ASN:HD22	2.07	0.58
1:A:250:LEU:HB3	1:A:273:VAL:HG13	1.86	0.58
1:B:338:VAL:HG21	1:B:354:ILE:HD11	1.86	0.58
1:B:40:TYR:CZ	1:B:326:MET:HA	2.38	0.58
1:A:340:THR:O	1:A:344:LYS:HB2	2.03	0.58
1:A:57:GLU:OE2	1:A:301:HIS:HE1	1.85	0.58
1:B:334:ARG:HG3	1:B:353:ILE:HG22	1.86	0.57
1:A:211:LEU:HD22	1:A:215:LYS:HD3	1.85	0.57
1:B:323:LEU:HD12	1:B:326:MET:CE	2.34	0.57
1:A:158:TYR:CD1	1:A:173:LEU:CD2	2.88	0.57
1:A:336:LEU:HG	1:A:397:MET:HG2	1.85	0.57
1:B:24:PHE:CE2	1:B:32:LYS:HE3	2.39	0.57
1:B:100:ALA:O	1:B:101:ARG:HD2	2.03	0.57
1:B:129:ARG:HD2	1:B:134:TRP:NE1	2.18	0.57
1:B:156:ARG:HD3	3:B:439:HOH:O	2.04	0.57
1:B:107:GLY:HA3	1:B:266:ARG:HH12	1.67	0.57
1:A:258:LYS:HG2	1:A:359:MET:HE3	1.87	0.56
1:B:356:GLN:OE1	1:B:361:SER:HA	2.05	0.56
1:B:258:LYS:NZ	2:B:410:PLA:HB22	2.20	0.56
1:A:373:LEU:HA	1:A:377:PHE:HD2	1.71	0.56
1:A:337:PHE:O	1:A:341:LEU:HB2	2.05	0.56
1:B:203:GLU:HG2	3:B:428:HOH:O	2.05	0.56
1:B:389:VAL:O	1:B:392:MET:HB2	2.05	0.56
1:A:159:ALA:HB3	1:A:172:ALA:O	2.06	0.56
1:A:338:VAL:HG21	1:A:354:ILE:HD11	1.88	0.56
1:A:370:VAL:HG21	1:A:383:ALA:HA	1.88	0.55
1:A:361:SER:OG	1:A:387:VAL:HG23	2.06	0.55
1:A:366:THR:HG22	1:A:369:GLN:N	2.20	0.55
1:B:321:GLN:HG3	3:B:430:HOH:O	2.06	0.55
1:A:134:TRP:CH2	1:A:179:GLU:HG2	2.41	0.55
1:B:187:LEU:CD1	1:B:220:LEU:HG	2.37	0.55
1:A:215:LYS:HG2	3:A:510:HOH:O	2.05	0.55
1:A:56:ALA:HB1	1:A:308:THR:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:TYR:O	1:B:168:LEU:HD23	2.07	0.55
1:B:24:PHE:CE1	1:B:34:ASN:HB2	2.42	0.54
1:B:40:TYR:CZ	1:B:326:MET:HG2	2.42	0.54
1:A:18:LEU:HD23	1:B:73:ILE:HD11	1.89	0.54
1:A:86:LEU:HD12	1:A:256:TYR:HE1	1.72	0.54
1:B:143:HIS:O	1:B:147:PHE:CD2	2.60	0.54
1:B:152:LEU:HD13	1:B:152:LEU:H	1.73	0.54
1:B:194:ASN:HD22	2:B:410:PLA:H2A1	1.72	0.54
1:A:143:HIS:O	1:A:147:PHE:CD2	2.61	0.54
1:A:21:ALA:HB1	3:A:509:HOH:O	2.08	0.54
1:A:292:ARG:HD2	1:A:292:ARG:O	2.07	0.54
1:A:366:THR:HB	1:A:369:GLN:CD	2.28	0.53
2:A:410:PLA:N	2:A:410:PLA:O3	2.40	0.53
1:A:125:VAL:HG11	1:A:185:VAL:HG23	1.90	0.53
1:A:266:ARG:HH22	2:A:410:PLA:P	2.30	0.53
1:B:165:ASN:HD21	1:B:167:THR:CG2	2.20	0.53
1:B:206:GLN:O	1:B:209:ALA:HB3	2.08	0.53
1:A:258:LYS:HG2	1:A:359:MET:CE	2.39	0.53
1:A:286:GLN:CD	1:B:12:ALA:HB2	2.28	0.53
1:A:129:ARG:HH22	1:A:181:GLN:HG3	1.71	0.53
1:A:258:LYS:HD3	1:A:263:TYR:CE1	2.43	0.53
1:B:123:THR:HG21	1:B:125:VAL:HG13	1.86	0.53
1:B:370:VAL:HG21	1:B:383:ALA:HA	1.91	0.52
1:B:332:ARG:CZ	1:B:332:ARG:HB2	2.39	0.52
1:A:189:HIS:CD2	1:A:194:ASN:H	2.27	0.52
1:A:129:ARG:HE	1:A:134:TRP:HE1	1.58	0.52
1:B:291:ILE:HG23	1:B:295:TYR:CZ	2.45	0.52
1:B:323:LEU:HA	1:B:326:MET:HE2	1.91	0.52
1:B:86:LEU:HD12	1:B:256:TYR:HE1	1.73	0.52
1:B:258:LYS:HZ3	2:B:410:PLA:H4A2	1.72	0.52
1:B:344:LYS:HD3	1:B:344:LYS:O	2.09	0.52
1:A:334:ARG:HD3	1:A:353:ILE:O	2.11	0.51
1:A:59:TYR:O	1:A:63:ASN:ND2	2.42	0.51
1:B:187:LEU:HD12	1:B:220:LEU:HG	1.91	0.51
1:A:195:PRO:HB3	1:A:386:ARG:HG3	1.91	0.51
1:A:9:ILE:CG2	1:B:122:ASN:HB3	2.40	0.51
1:B:237:ALA:HB3	1:B:241:ARG:HH21	1.75	0.51
1:B:29:ARG:HD2	3:B:414:HOH:O	2.10	0.51
1:B:8:ASN:HB3	3:B:461:HOH:O	2.10	0.51
1:A:73:ILE:HG21	1:A:288:LYS:HG2	1.92	0.51
1:A:319:TRP:CZ2	1:A:323:LEU:HD22	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:PRO:HD2	1:A:295:TYR:CZ	2.46	0.51
1:A:5:MET:N	1:A:7:GLU:OE1	2.44	0.51
1:B:316:ARG:HD2	1:B:320:GLU:OE2	2.11	0.51
1:B:102:THR:CG2	1:B:269:ALA:HB1	2.42	0.50
1:B:37:ILE:O	1:B:388:ASN:ND2	2.44	0.50
1:A:168:LEU:HB3	1:A:198:ILE:HG21	1.92	0.50
1:A:369:GLN:HG2	1:A:408:VAL:O	2.11	0.50
1:B:71:LEU:HB2	3:B:411:HOH:O	2.11	0.50
1:A:194:ASN:HD22	2:A:410:PLA:H2A1	1.77	0.50
1:A:314:ALA:O	1:A:318:ILE:HG12	2.12	0.50
1:B:393:THR:HG22	1:B:396:ASN:H	1.76	0.50
1:B:36:GLY:HA3	3:B:511:HOH:O	2.10	0.50
1:B:229:ALA:O	1:B:236:ASP:OD1	2.30	0.50
1:A:53:VAL:O	1:A:57:GLU:HG3	2.11	0.50
1:B:123:THR:CG2	1:B:125:VAL:H	2.24	0.50
1:A:133:VAL:HG22	1:A:185:VAL:HB	1.94	0.49
1:B:158:TYR:HD1	1:B:173:LEU:HD22	1.77	0.49
1:B:79:PHE:CD2	1:B:104:GLN:HG3	2.47	0.49
1:A:334:ARG:NH2	1:A:358:GLY:O	2.45	0.49
1:A:398:ALA:N	1:A:399:PRO:HD2	2.26	0.49
1:B:123:THR:HG23	1:B:124:SER:N	2.27	0.49
1:B:200:PRO:HG2	1:B:205:TRP:CE2	2.47	0.49
1:B:277:SER:O	1:B:280:VAL:HG12	2.13	0.49
1:B:179:GLU:O	1:B:179:GLU:HG3	2.11	0.49
1:B:41:LYS:O	1:B:329:ARG:NH1	2.45	0.49
1:A:67:THR:OG1	1:B:47:THR:O	2.22	0.49
1:B:80:GLY:O	1:B:83:THR:HB	2.12	0.49
1:A:44:THR:OG1	1:A:46:LYS:HG3	2.13	0.49
1:B:223:PHE:CD2	1:B:240:LEU:HD12	2.47	0.49
1:B:212:SER:HG	1:B:247:HIS:CE1	2.31	0.48
1:A:139:SER:N	3:A:477:HOH:O	2.38	0.48
1:A:61:LEU:CD1	1:B:61:LEU:HD12	2.34	0.48
1:A:97:ASP:O	1:A:98:LYS:HB2	2.14	0.48
1:B:129:ARG:CZ	1:B:129:ARG:HB3	2.43	0.48
1:A:205:TRP:HE3	1:A:243:PHE:HE1	1.61	0.48
1:A:260:PHE:HB3	1:A:262:LEU:CD2	2.29	0.48
1:B:348:ARG:NH2	1:B:369:GLN:HE22	2.11	0.48
1:B:372:ARG:HG2	1:B:408:VAL:CG1	2.42	0.48
1:A:188:PHE:CE1	1:A:208:LEU:HD13	2.49	0.48
1:B:393:THR:CG2	1:B:395:ASP:H	2.27	0.48
1:A:231:GLY:O	1:A:236:ASP:HB2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:ASN:OD1	1:A:315:LEU:HB2	2.14	0.47
1:A:350:PHE:HB3	1:A:353:ILE:CD1	2.42	0.47
1:B:323:LEU:HA	1:B:326:MET:HE3	1.96	0.47
1:B:229:ALA:HB2	1:B:357:ASN:HB3	1.95	0.47
1:A:366:THR:HB	1:A:369:GLN:OE1	2.14	0.47
1:B:5:MET:N	1:B:7:GLU:OE1	2.47	0.47
1:A:144:LYS:HG2	1:A:155:VAL:HG21	1.96	0.47
1:A:362:PHE:HE1	1:A:384:SER:HB2	1.79	0.47
1:A:258:LYS:HD3	1:A:263:TYR:HE1	1.78	0.47
1:A:122:ASN:CG	1:A:286:GLN:HG2	2.35	0.47
1:B:29:ARG:HB3	1:B:30:PRO:HD2	1.96	0.47
1:A:222:ASP:OD2	2:A:410:PLA:N1	2.48	0.47
1:A:291:ILE:HG23	1:A:295:TYR:CZ	2.50	0.47
1:B:272:LEU:HD12	1:B:287:MET:SD	2.55	0.47
1:B:276:ASP:O	1:B:280:VAL:HG12	2.15	0.47
1:B:362:PHE:HA	1:B:362:PHE:HD1	1.62	0.47
1:B:270:CYS:SG	1:B:287:MET:CE	3.03	0.47
1:A:359:MET:N	3:A:465:HOH:O	2.41	0.47
1:A:93:ALA:O	1:A:97:ASP:HB2	2.15	0.47
1:B:120:ALA:O	1:B:121:LYS:HB2	2.15	0.47
1:A:104:GLN:NE2	3:A:495:HOH:O	2.43	0.46
1:B:133:VAL:CB	1:B:185:VAL:HG22	2.43	0.46
1:B:5:MET:C	1:B:8:ASN:HD21	2.18	0.46
1:B:15:ASP:CG	1:B:16:PRO:HD2	2.35	0.46
1:B:372:ARG:CZ	1:B:408:VAL:HG12	2.46	0.46
1:A:403:ALA:O	1:A:404:ILE:C	2.54	0.46
1:B:133:VAL:HA	1:B:185:VAL:O	2.15	0.46
1:B:335:GLN:O	1:B:338:VAL:HG22	2.16	0.46
1:B:372:ARG:HG2	1:B:408:VAL:HG11	1.98	0.46
1:B:401:CYS:O	1:B:405:VAL:HB	2.16	0.46
1:A:302:GLY:O	1:A:306:VAL:HG23	2.15	0.46
1:B:355:LYS:HD3	1:B:355:LYS:C	2.36	0.46
1:A:20:LEU:HD23	1:A:20:LEU:HA	1.67	0.45
1:B:229:ALA:O	1:B:230:ARG:HB2	2.16	0.45
1:B:340:THR:O	1:B:344:LYS:HB2	2.16	0.45
1:B:79:PHE:CE2	1:B:104:GLN:HG3	2.51	0.45
1:B:144:LYS:HG2	1:B:144:LYS:HZ2	1.57	0.45
1:B:333:MET:HE2	1:B:336:LEU:HD23	1.97	0.45
1:B:370:VAL:O	1:B:370:VAL:HG23	2.16	0.45
1:B:266:ARG:NH2	2:B:410:PLA:O1P	2.45	0.45
1:B:137:ASN:HD22	1:B:138:PRO:HA	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:ASP:O	1:A:32:LYS:NZ	2.38	0.45
1:B:192:CYS:SG	1:B:229:ALA:HB2	2.57	0.45
1:B:338:VAL:CG1	1:B:354:ILE:HD11	2.45	0.45
1:A:68:LYS:HB2	3:B:424:HOH:O	2.16	0.45
1:A:137:ASN:O	1:A:160:TYR:HB3	2.17	0.45
1:A:35:LEU:HD11	1:A:400:LEU:HD13	1.99	0.45
1:B:115:ALA:O	1:B:118:PHE:HB3	2.17	0.45
1:B:373:LEU:HA	1:B:373:LEU:HD12	1.71	0.45
1:B:237:ALA:HB3	1:B:241:ARG:HE	1.82	0.45
1:A:158:TYR:HD1	1:A:173:LEU:CD2	2.30	0.44
1:A:372:ARG:O	1:A:376:GLU:HB2	2.16	0.44
1:A:338:VAL:HG21	1:A:354:ILE:CD1	2.47	0.44
1:A:373:LEU:HD23	1:A:377:PHE:CD2	2.52	0.44
1:A:377:PHE:HE2	1:A:408:VAL:CG1	2.30	0.44
1:B:40:TYR:CG	1:B:41:LYS:N	2.85	0.44
1:A:24:PHE:CZ	1:A:34:ASN:HB2	2.51	0.44
1:A:41:LYS:O	1:A:329:ARG:NH1	2.49	0.44
1:A:61:LEU:O	1:B:54:LYS:NZ	2.49	0.44
1:B:123:THR:CG2	1:B:124:SER:N	2.80	0.44
1:B:190:GLY:HA2	1:B:221:PHE:CE1	2.53	0.44
1:B:373:LEU:HG	1:B:379:VAL:HB	1.99	0.44
1:A:220:LEU:HD11	1:A:251:ILE:HG21	2.00	0.44
1:A:305:VAL:O	1:A:308:THR:HB	2.17	0.44
1:B:113:ARG:O	1:B:116:ALA:HB3	2.17	0.44
1:B:119:LEU:O	1:B:123:THR:HB	2.18	0.44
1:B:314:ALA:O	1:B:317:ALA:HB3	2.17	0.44
1:B:133:VAL:O	1:B:155:VAL:HA	2.17	0.44
1:A:305:VAL:HG12	1:A:306:VAL:N	2.32	0.44
1:A:334:ARG:NH1	3:A:513:HOH:O	2.51	0.44
1:A:118:PHE:CE1	1:A:119:LEU:HD23	2.53	0.44
1:B:237:ALA:CB	1:B:241:ARG:HE	2.30	0.43
1:A:348:ARG:HH22	1:A:369:GLN:HE22	1.66	0.43
1:B:348:ARG:HH22	1:B:369:GLN:HE22	1.65	0.43
1:B:121:LYS:HD3	1:B:121:LYS:C	2.39	0.43
1:B:41:LYS:HG2	1:B:45:GLY:HA2	2.00	0.43
1:B:158:TYR:CD1	1:B:173:LEU:HD22	2.53	0.43
1:A:397:MET:HG3	1:A:397:MET:O	2.18	0.43
1:A:76:ILE:HA	1:A:77:PRO:HD3	1.89	0.43
1:A:109:THR:O	1:B:294:ASN:O	2.37	0.43
1:A:126:LYS:HA	1:A:126:LYS:HD2	1.69	0.43
1:A:100:ALA:C	1:A:101:ARG:HG2	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:GLU:H	1:B:78:GLU:CD	2.22	0.43
1:B:39:VAL:HG23	1:B:263:TYR:CE1	2.54	0.43
1:B:42:ASP:HB2	1:B:43:GLU:OE2	2.18	0.43
1:A:34:ASN:HA	1:A:380:TYR:HB2	2.01	0.43
1:B:53:VAL:HG21	3:B:443:HOH:O	2.18	0.43
1:A:350:PHE:HZ	1:A:409:LEU:HD21	1.83	0.42
1:B:229:ALA:CB	1:B:357:ASN:HB3	2.49	0.42
1:A:113:ARG:O	1:A:113:ARG:HG3	2.18	0.42
1:A:308:THR:HG22	1:A:309:ILE:N	2.33	0.42
1:B:270:CYS:SG	1:B:287:MET:HE2	2.59	0.42
1:B:101:ARG:NH2	1:B:281:ASP:OD1	2.53	0.42
1:B:270:CYS:SG	1:B:287:MET:HE1	2.59	0.42
1:B:292:ARG:HG3	1:B:292:ARG:O	2.19	0.42
1:B:324:THR:O	1:B:324:THR:HG22	2.18	0.42
1:B:354:ILE:HG22	1:B:354:ILE:O	2.19	0.42
1:B:372:ARG:NH2	1:B:407:ALA:O	2.53	0.42
1:A:202:LEU:HD22	1:A:206:GLN:NE2	2.35	0.42
1:A:223:PHE:HD1	1:A:223:PHE:HA	1.73	0.42
1:A:35:LEU:HD11	1:A:400:LEU:CD1	2.50	0.42
1:B:240:LEU:O	1:B:240:LEU:HD23	2.20	0.42
1:B:30:PRO:HG2	1:B:31:GLY:H	1.84	0.42
1:B:6:PHE:HD1	1:B:9:ILE:HD11	1.84	0.42
1:B:348:ARG:O	1:B:350:PHE:CD1	2.72	0.42
1:B:152:LEU:HD13	1:B:152:LEU:N	2.35	0.42
1:A:344:LYS:HA	1:A:344:LYS:HD2	1.86	0.42
1:B:195:PRO:HB3	1:B:386:ARG:CG	2.48	0.42
1:A:24:PHE:CD1	1:A:34:ASN:HB2	2.54	0.42
1:A:256:TYR:HB2	1:A:267:VAL:HB	2.03	0.41
1:A:67:THR:OG1	1:A:68:LYS:N	2.54	0.41
1:A:5:MET:HE2	1:B:123:THR:HG23	2.02	0.41
1:A:109:THR:O	1:A:109:THR:CG2	2.67	0.41
1:A:373:LEU:HA	1:A:373:LEU:HD23	1.80	0.41
1:A:9:ILE:HG23	1:B:122:ASN:HB3	2.00	0.41
1:B:112:LEU:HB3	1:B:146:VAL:HG21	2.02	0.41
1:B:341:LEU:HD12	1:B:405:VAL:HG23	2.02	0.41
1:B:123:THR:HG23	1:B:125:VAL:HG13	1.93	0.41
1:B:392:MET:HG2	1:B:400:LEU:CD2	2.50	0.41
1:A:133:VAL:O	1:A:155:VAL:HA	2.21	0.41
1:A:264:ASN:HB2	1:B:71:LEU:HG	2.02	0.41
1:B:6:PHE:HA	1:B:9:ILE:HG13	2.02	0.41
1:A:211:LEU:HD23	1:A:211:LEU:HA	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:LEU:HD12	1:A:287:MET:SD	2.61	0.41
1:B:393:THR:HG23	1:B:395:ASP:H	1.85	0.41
1:B:5:MET:HE2	1:B:5:MET:HB2	1.72	0.41
1:B:338:VAL:CG2	1:B:339:ASN:N	2.84	0.41
1:B:225:TYR:OH	2:B:410:PLA:HB23	2.21	0.41
1:B:260:PHE:CD2	1:B:306:VAL:HG22	2.56	0.41
1:B:359:MET:HB3	1:B:388:ASN:OD1	2.21	0.41
1:B:383:ALA:HB2	3:B:467:HOH:O	2.21	0.40
1:A:205:TRP:HE3	1:A:243:PHE:CE1	2.39	0.40
1:B:209:ALA:O	1:B:213:VAL:CG2	2.62	0.40
1:B:280:VAL:HG13	1:B:281:ASP:N	2.35	0.40
1:A:18:LEU:HD13	1:B:292:ARG:NH2	2.36	0.40
1:A:39:VAL:HG23	1:A:263:TYR:CE1	2.57	0.40
1:A:301:HIS:O	1:A:304:SER:HB2	2.21	0.40
1:A:30:PRO:HG2	1:A:31:GLY:H	1.86	0.40
1:B:161:TYR:CE2	1:B:362:PHE:HD2	2.39	0.40
1:B:301:HIS:HB3	1:B:302:GLY:H	1.78	0.40
1:B:312:ASN:OD1	1:B:315:LEU:HG	2.22	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	394/396 (100%)	345 (88%)	37 (9%)	12 (3%)	4	7
1	B	394/396 (100%)	335 (85%)	44 (11%)	15 (4%)	3	4
All	All	788/792 (100%)	680 (86%)	81 (10%)	27 (3%)	3	5

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	92	SER
1	B	136	SER
1	B	152	LEU
1	B	301	HIS
1	A	36	GLY
1	A	50	LEU
1	A	121	LYS
1	A	301	HIS
1	B	26	ALA
1	B	30	PRO
1	B	121	LYS
1	B	390	ALA
1	A	141	PRO
1	A	215	LYS
1	A	338	VAL
1	A	378	GLY
1	B	36	GLY
1	B	37	ILE
1	B	43	GLU
1	B	70	TYR
1	B	141	PRO
1	A	43	GLU
1	A	201	THR
1	B	90	LYS
1	A	404	ILE
1	B	195	PRO
1	A	195	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	320/320 (100%)	257 (80%)	63 (20%)	1	2
1	B	320/320 (100%)	253 (79%)	67 (21%)	1	2
All	All	640/640 (100%)	510 (80%)	130 (20%)	1	2

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	MET
1	A	8	ASN
1	A	9	ILE
1	A	23	LEU
1	A	28	GLU
1	A	39	VAL
1	A	41	LYS
1	A	44	THR
1	A	61	LEU
1	A	62	GLU
1	A	64	GLU
1	A	87	LEU
1	A	90	LYS
1	A	96	ASN
1	A	98	LYS
1	A	104	GLN
1	A	118	PHE
1	A	121	LYS
1	A	124	SER
1	A	144	LYS
1	A	154	GLU
1	A	155	VAL
1	A	165	ASN
1	A	168	LEU
1	A	170	PHE
1	A	174	ILE
1	A	179	GLU
1	A	206	GLN
1	A	211	LEU
1	A	215	LYS
1	A	218	LEU
1	A	223	PHE
1	A	240	LEU
1	A	250	LEU
1	A	258	LYS
1	A	259	ASN
1	A	262	LEU
1	A	273	VAL
1	A	292	ARG
1	A	296	SER
1	A	297	ASN
1	A	313	ASP
1	A	316	ARG

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Mol	Chain	Res	Type
1	A	329	ARG
1	A	332	ARG
1	A	341	LEU
1	A	342	GLN
1	A	343	GLU
1	A	347	ASN
1	A	348	ARG
1	A	353	ILE
1	A	360	PHE
1	A	362	PHE
1	A	367	LYS
1	A	370	VAL
1	A	371	LEU
1	A	372	ARG
1	A	374	ARG
1	A	375	GLU
1	A	376	GLU
1	A	384	SER
1	A	387	VAL
1	A	388	ASN
1	B	5	MET
1	B	8	ASN
1	B	10	THR
1	B	28	GLU
1	B	29	ARG
1	B	41	LYS
1	B	55	LYS
1	B	60	LEU
1	B	61	LEU
1	B	67	THR
1	B	86	LEU
1	B	87	LEU
1	B	90	LYS
1	B	98	LYS
1	B	112	LEU
1	B	114	VAL
1	B	119	LEU
1	B	121	LYS
1	B	123	THR
1	B	129	ARG
1	B	133	VAL
1	B	144	LYS

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Mol	Chain	Res	Type
1	B	146	VAL
1	B	152	LEU
1	B	165	ASN
1	B	168	LEU
1	B	173	LEU
1	B	179	GLU
1	B	185	VAL
1	B	210	GLN
1	B	215	LYS
1	B	223	PHE
1	B	233	LEU
1	B	234	GLU
1	B	235	GLU
1	B	240	LEU
1	B	246	MET
1	B	254	SER
1	B	259	ASN
1	B	262	LEU
1	B	266	ARG
1	B	273	VAL
1	B	277	SER
1	B	294	ASN
1	B	297	ASN
1	B	304	SER
1	B	311	SER
1	B	313	ASP
1	B	316	ARG
1	B	329	ARG
1	B	332	ARG
1	B	333	MET
1	B	334	ARG
1	B	341	LEU
1	B	343	GLU
1	B	344	LYS
1	B	347	ASN
1	B	348	ARG
1	B	349	ASP
1	B	355	LYS
1	B	362	PHE
1	B	370	VAL
1	B	371	LEU
1	B	372	ARG

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Mol	Chain	Res	Type
1	B	373	LEU
1	B	384	SER
1	B	393	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	58	GLN
1	A	63	ASN
1	A	104	GLN
1	A	137	ASN
1	A	175	ASN
1	A	181	GLN
1	A	189	HIS
1	A	226	GLN
1	A	247	HIS
1	A	259	ASN
1	A	301	HIS
1	A	339	ASN
1	A	357	ASN
1	A	388	ASN
1	B	8	ASN
1	B	69	ASN
1	B	137	ASN
1	B	165	ASN
1	B	175	ASN
1	B	226	GLN
1	B	259	ASN
1	B	294	ASN
1	B	331	GLN

### 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

2 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
2	PLA	B	410	-	18,25,25	2.29	4 (22%)	24,37,37	3.19	12 (50%)
2	PLA	A	410	-	18,25,25	2.15	4 (22%)	24,37,37	3.31	14 (58%)

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
2	PLA	B	410	-	-	6/15/23/23	0/1/1/1
2	PLA	A	410	-	-	7/15/23/23	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	410	PLA	C4A-C4	-6.76	1.43	1.51
2	A	410	PLA	C4A-C4	-5.87	1.44	1.51
2	B	410	PLA	C3-C4	-3.84	1.34	1.40
2	A	410	PLA	CB1-CA	-3.55	1.50	1.55
2	B	410	PLA	CB1-CA	-3.45	1.50	1.55
2	A	410	PLA	CB2-CA	-3.16	1.51	1.53
2	A	410	PLA	C3-C4	-2.62	1.36	1.40
2	B	410	PLA	C5-C4	2.37	1.43	1.40

All (26) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	410	PLA	C4A-C4-C5	6.36	126.78	119.71
2	A	410	PLA	C2A-C2-C3	5.43	127.59	120.89
2	B	410	PLA	C2A-C2-C3	5.40	127.56	120.89
2	A	410	PLA	C4A-C4-C5	5.40	125.71	119.71
2	B	410	PLA	C4A-C4-C3	-5.36	114.31	120.04
2	A	410	PLA	C6-C5-C4	5.15	121.77	118.12
2	A	410	PLA	CB2-CA-CB1	-5.01	104.00	110.84
2	B	410	PLA	O2P-P-O4P	4.91	119.80	106.73
2	A	410	PLA	O2P-P-O4P	4.85	119.65	106.73
2	A	410	PLA	O4P-C5A-C5	4.57	118.05	109.35
2	B	410	PLA	C3-C2-N1	-4.39	115.09	120.77
2	A	410	PLA	C3-C2-N1	-4.36	115.14	120.77
2	B	410	PLA	C5-C6-N1	-4.30	116.65	123.82
2	A	410	PLA	O3P-P-O1P	4.27	127.40	110.68
2	B	410	PLA	C6-C5-C4	4.21	121.09	118.12
2	A	410	PLA	C5-C6-N1	-4.15	116.91	123.82
2	B	410	PLA	C6-N1-C2	4.05	126.67	119.17
2	A	410	PLA	C6-N1-C2	3.98	126.55	119.17
2	B	410	PLA	O3P-P-O1P	3.83	125.66	110.68
2	B	410	PLA	CA-CB1-CG	-2.99	109.01	114.94
2	A	410	PLA	O4P-P-O1P	-2.90	98.35	106.47
2	B	410	PLA	CB2-CA-CB1	2.83	114.71	110.84
2	A	410	PLA	C4A-C4-C3	-2.71	117.14	120.04
2	A	410	PLA	C5A-C5-C6	-2.28	115.62	119.37
2	B	410	PLA	O3P-P-O4P	-2.17	100.96	106.73
2	A	410	PLA	C3-C4-C5	-2.04	116.77	118.72

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	410	PLA	C4-C5-C5A-O4P
2	B	410	PLA	C6-C5-C5A-O4P
2	A	410	PLA	C4-C5-C5A-O4P
2	A	410	PLA	C6-C5-C5A-O4P
2	B	410	PLA	C5-C4-C4A-N
2	A	410	PLA	C5-C4-C4A-N
2	B	410	PLA	C5A-O4P-P-O1P
2	A	410	PLA	C5A-O4P-P-O1P
2	B	410	PLA	C5A-O4P-P-O2P
2	A	410	PLA	CB2-CA-N-C4A
2	B	410	PLA	C3-C4-C4A-N
2	A	410	PLA	C3-C4-C4A-N

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Mol	Chain	Res	Type	Atoms
2	A	410	PLA	C5A-O4P-P-O2P

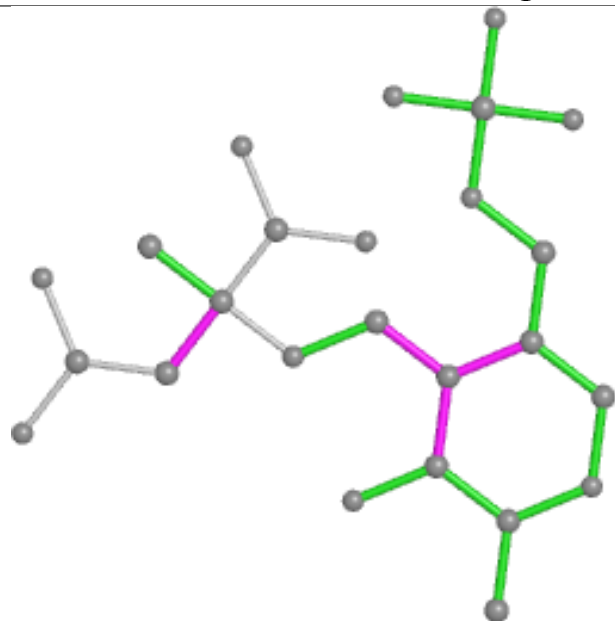
There are no ring outliers.

2 monomers are involved in 15 short contacts:

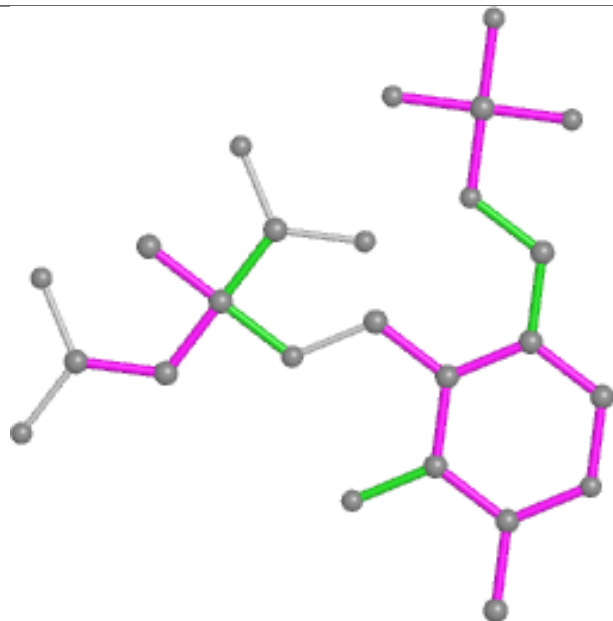
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	410	PLA	10	0
2	A	410	PLA	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

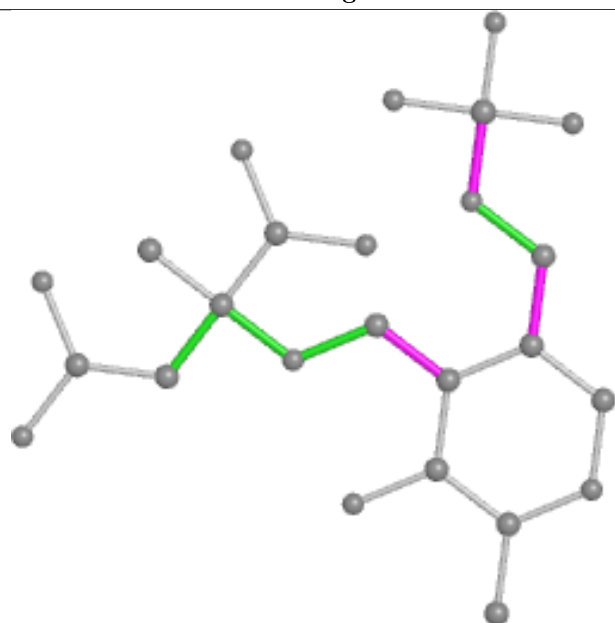
## Ligand PLA B 410



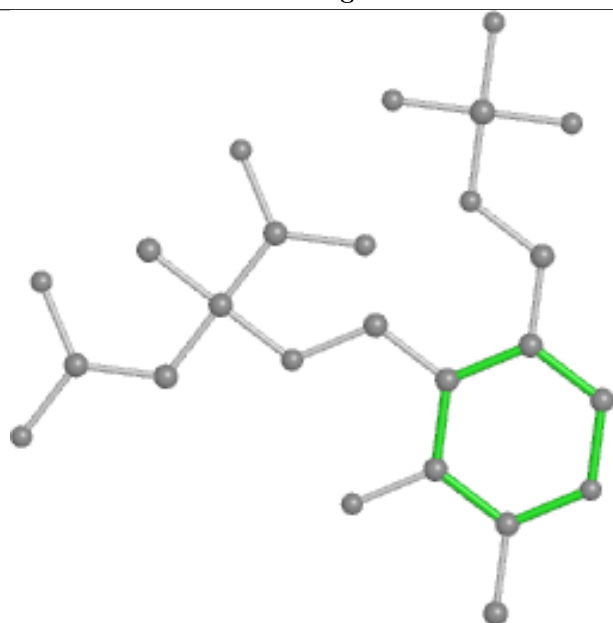
Bond lengths



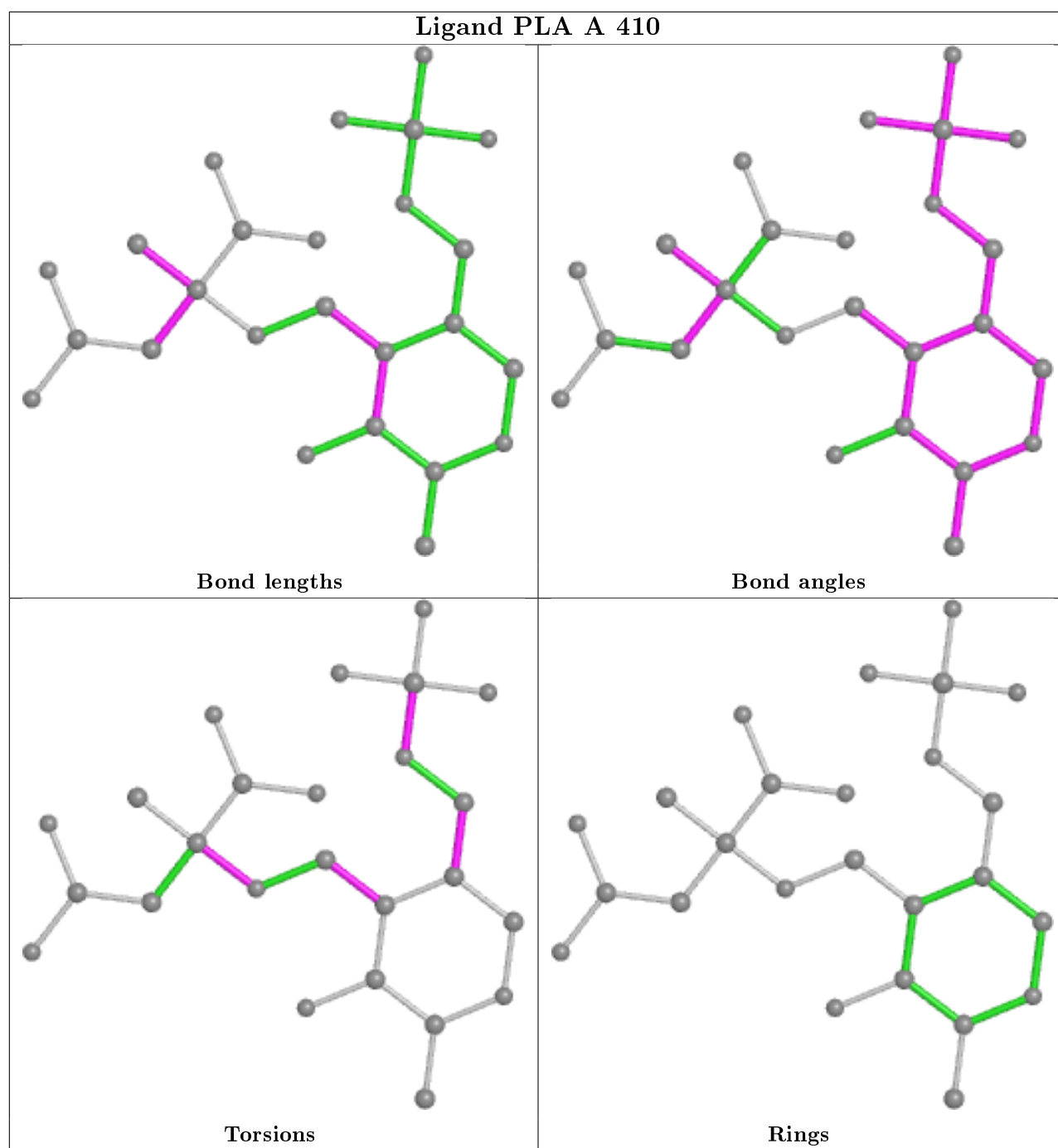
Bond angles



Torsions



Rings



## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	396/396 (100%)	-1.19	0 100 100	8, 30, 67, 86	0
1	B	396/396 (100%)	-1.17	0 100 100	4, 31, 67, 81	0
All	All	792/792 (100%)	-1.18	0 100 100	4, 31, 67, 86	0

There are no RSRZ outliers to report.

### 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](#)

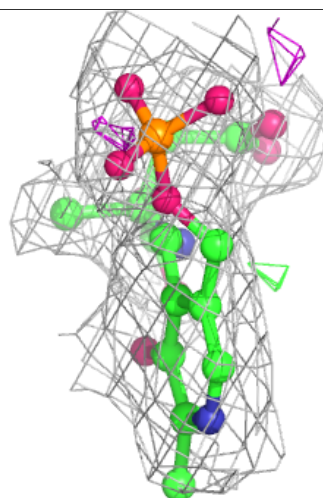
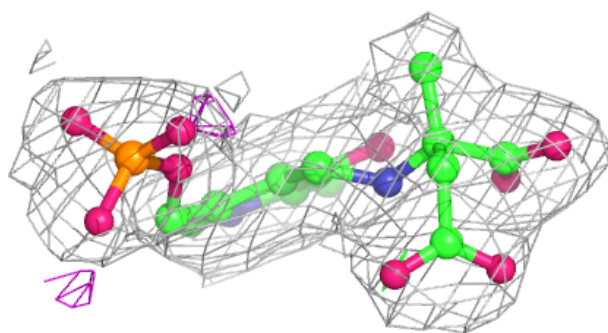
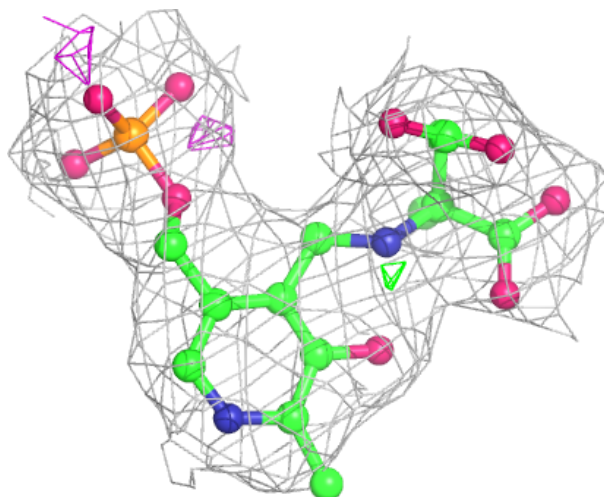
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

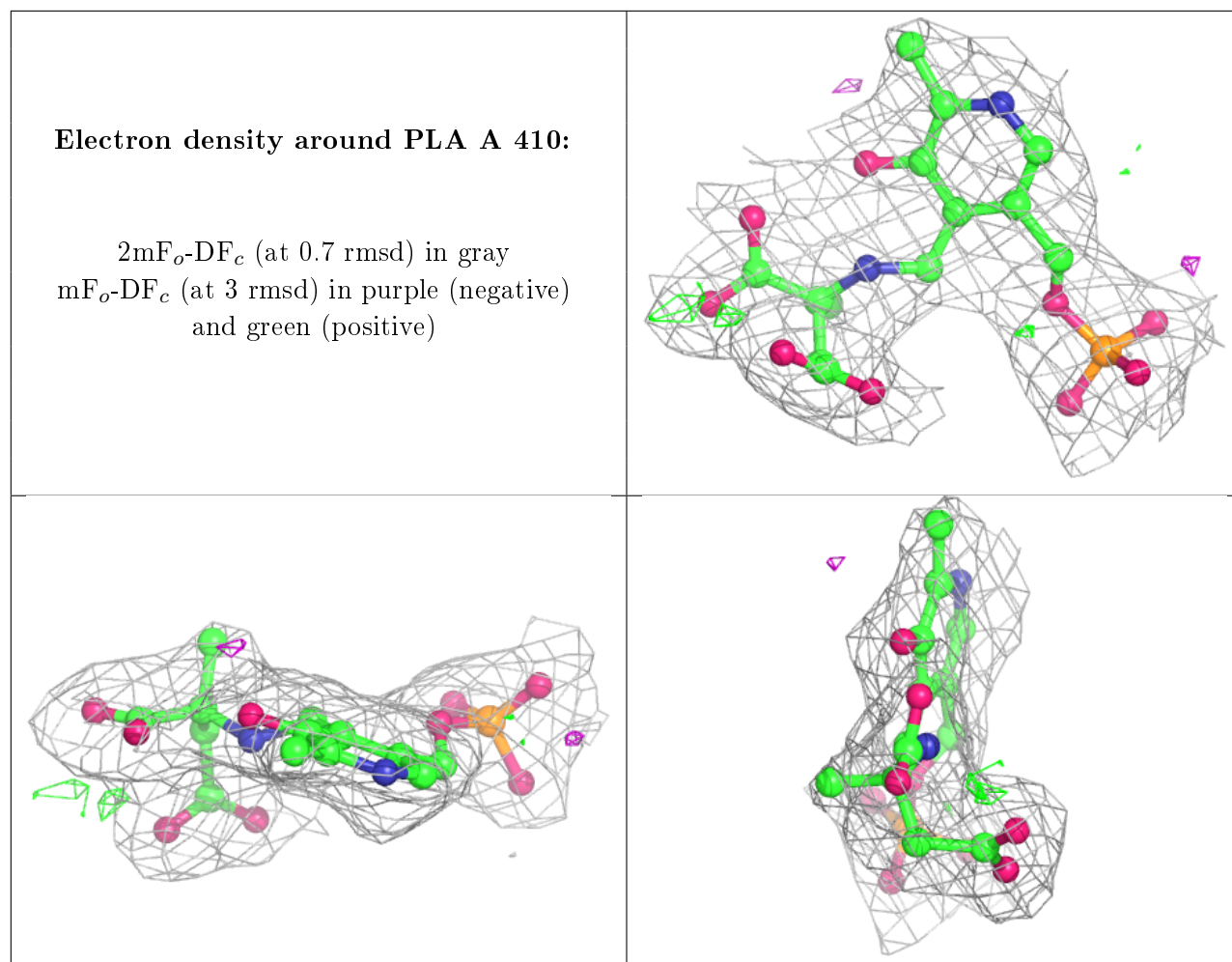
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	PLA	B	410	25/25	0.98	0.11	9,20,48,76	0
2	PLA	A	410	25/25	0.98	0.10	7,22,40,49	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around PLA B 410:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers ⓘ

There are no such residues in this entry.