Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 05:18 PM EDT

PDB ID	:	1TDJ
Title	:	THREONINE DEAMINASE (BIOSYNTHETIC) FROM E. COLI
Authors	:	Gallagher, D.T.; Gilliland, G.L.; Xiao, G.; Eisenstein, E.
Deposited on	:	1998-03-27
Resolution	:	2.80 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

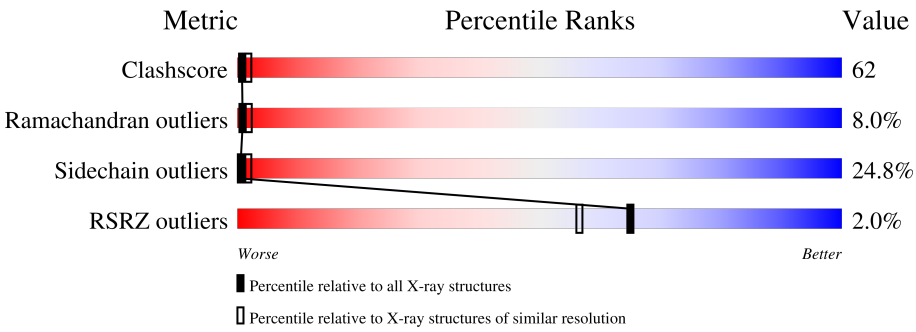
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

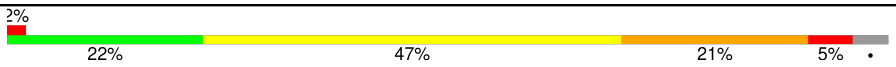
The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 of the lower bar indicate the fraction of residues that contain outliers for >=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 <=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	514	

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	PLP	A	962	-	X	-	-

2 Entry composition (i)

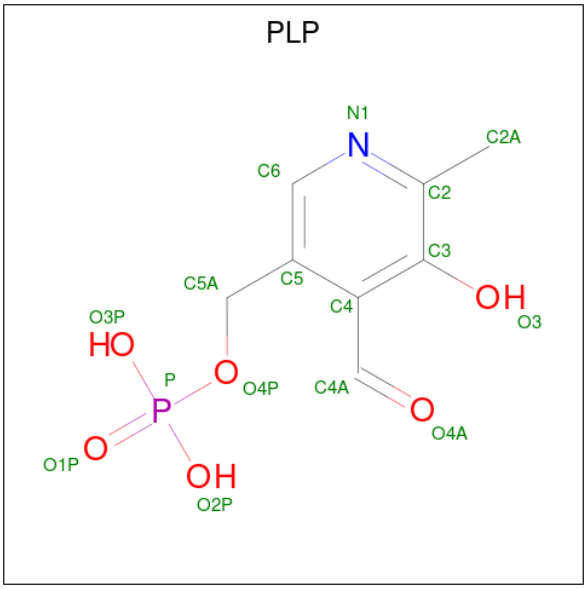
There are 3 unique types of molecules in this entry. The entry contains 3848 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BIOSYNTHETIC THREONINE DEAMINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	494	3800	2412	669	701	18	0	0	0

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	15	8	1	5	1	0	0

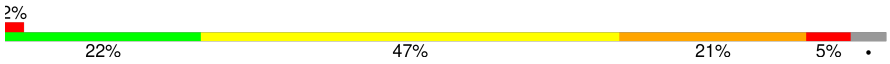
- Molecule 3 is water.

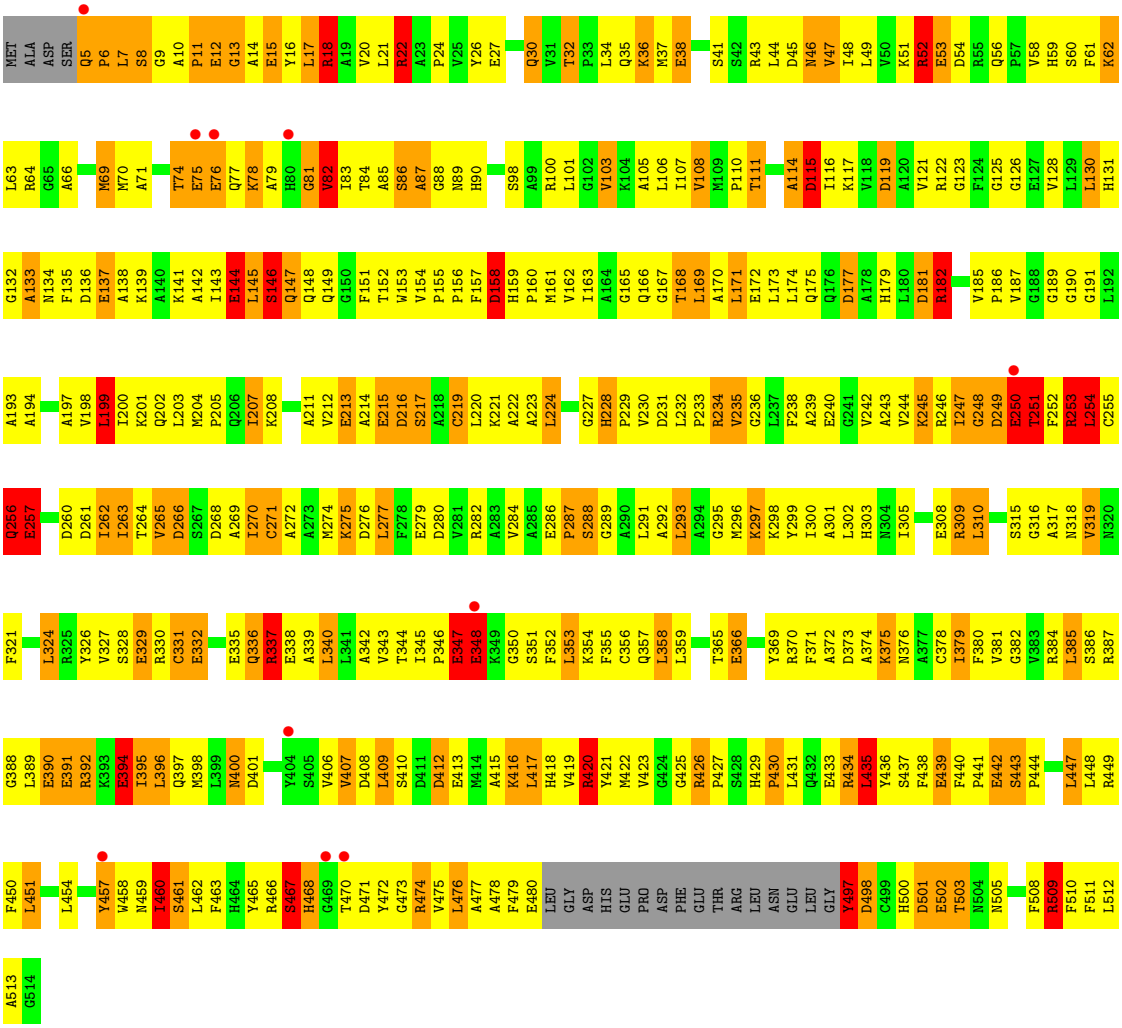
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	33	33	33	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: BIOSYNTHETIC THREONINE DEAMINASE

Chain A: 



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	85.10Å 90.80Å 162.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	16.00 – 2.80 15.96 – 2.71	Depositor EDS
% Data completeness (in resolution range)	86.0 (16.00-2.80) 79.5 (15.96-2.71)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	0.71 (at 2.69Å)	Xtriage
Refinement program	TNT 5E	Depositor
R, R_{free}	0.200 , 0.340 0.192 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	36.7	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.21 , 93.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3848	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: PLP

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	1.29	28/3875 (0.7%)	1.71	85/5238 (1.6%)

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	A	2	1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	15	GLU	CD-OE1	7.75	1.34	1.25
1	A	53	GLU	CD-OE1	7.73	1.34	1.25
1	A	413	GLU	CD-OE2	7.55	1.33	1.25
1	A	76	GLU	CD-OE2	7.28	1.33	1.25
1	A	348	GLU	CD-OE2	7.26	1.33	1.25
1	A	390	GLU	CD-OE1	7.18	1.33	1.25
1	A	257	GLU	CD-OE1	7.08	1.33	1.25
1	A	250	GLU	CD-OE1	6.90	1.33	1.25
1	A	329	GLU	CD-OE1	-6.57	1.18	1.25
1	A	279	GLU	CD-OE1	6.49	1.32	1.25
1	A	240	GLU	CD-OE1	6.38	1.32	1.25
1	A	439	GLU	CD-OE2	6.18	1.32	1.25
1	A	213	GLU	CD-OE2	6.13	1.32	1.25
1	A	38	GLU	CD-OE2	6.06	1.32	1.25
1	A	12	GLU	CD-OE2	6.03	1.32	1.25
1	A	137	GLU	CD-OE1	5.94	1.32	1.25
1	A	75	GLU	CD-OE1	5.64	1.31	1.25
1	A	442	GLU	CD-OE2	5.49	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	366	GLU	CD-OE1	5.48	1.31	1.25
1	A	27	GLU	CD-OE1	-5.40	1.19	1.25
1	A	391	GLU	CD-OE1	5.37	1.31	1.25
1	A	215	GLU	CD-OE1	5.32	1.31	1.25
1	A	144	GLU	CD-OE2	5.31	1.31	1.25
1	A	332	GLU	CD-OE2	5.27	1.31	1.25
1	A	502	GLU	CD-OE2	5.21	1.31	1.25
1	A	45	ASP	CG-OD2	5.12	1.37	1.25
1	A	347	GLU	CD-OE2	5.07	1.31	1.25
1	A	146	SER	CB-OG	-5.03	1.35	1.42

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	253	ARG	NE-CZ-NH1	9.83	125.22	120.30
1	A	408	ASP	CB-CG-OD2	-9.26	109.97	118.30
1	A	158	ASP	CB-CG-OD1	8.79	126.21	118.30
1	A	45	ASP	CB-CG-OD2	-8.41	110.73	118.30
1	A	158	ASP	CB-CG-OD2	-8.33	110.80	118.30
1	A	426	ARG	NE-CZ-NH2	-8.29	116.16	120.30
1	A	249	ASP	CB-CG-OD2	-8.06	111.05	118.30
1	A	119	ASP	CB-CG-OD1	-8.05	111.05	118.30
1	A	280	ASP	CB-CG-OD1	-8.05	111.05	118.30
1	A	337	ARG	NE-CZ-NH2	8.04	124.32	120.30
1	A	52	ARG	NE-CZ-NH1	-8.03	116.28	120.30
1	A	373	ASP	CB-CG-OD1	-7.86	111.22	118.30
1	A	501	ASP	CB-CG-OD2	-7.80	111.28	118.30
1	A	471	ASP	CB-CG-OD1	7.55	125.09	118.30
1	A	280	ASP	N-CA-CB	7.53	124.15	110.60
1	A	276	ASP	CB-CG-OD2	-7.51	111.55	118.30
1	A	509	ARG	NE-CZ-NH1	7.43	124.01	120.30
1	A	18	ARG	NE-CZ-NH1	7.33	123.96	120.30
1	A	471	ASP	CB-CG-OD2	-7.30	111.73	118.30
1	A	412	ASP	CB-CG-OD2	-7.24	111.79	118.30
1	A	52	ARG	CD-NE-CZ	-7.21	113.50	123.60
1	A	373	ASP	CB-CG-OD2	7.12	124.71	118.30
1	A	254	LEU	N-CA-CB	7.11	124.61	110.40
1	A	234	ARG	NE-CZ-NH1	7.04	123.82	120.30
1	A	254	LEU	CB-CA-C	7.03	123.55	110.20
1	A	501	ASP	CB-CG-OD1	6.97	124.57	118.30
1	A	260	ASP	CB-CG-OD2	-6.96	112.04	118.30
1	A	216	ASP	CB-CG-OD2	-6.84	112.14	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	115	ASP	CB-CG-OD1	-6.82	112.16	118.30
1	A	216	ASP	CB-CG-OD1	6.80	124.42	118.30
1	A	498	ASP	CB-CG-OD1	-6.56	112.40	118.30
1	A	412	ASP	CB-CG-OD1	6.49	124.14	118.30
1	A	416	LYS	N-CA-CB	6.41	122.14	110.60
1	A	136	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	A	474	ARG	NE-CZ-NH2	6.37	123.48	120.30
1	A	119	ASP	CB-CG-OD2	6.33	124.00	118.30
1	A	146	SER	CB-CA-C	-6.24	98.24	110.10
1	A	249	ASP	CB-CG-OD1	6.24	123.92	118.30
1	A	347	GLU	N-CA-CB	6.21	121.77	110.60
1	A	136	ASP	CB-CG-OD1	6.17	123.85	118.30
1	A	251	THR	CB-CA-C	6.10	128.06	111.60
1	A	26	TYR	CB-CG-CD1	-6.08	117.35	121.00
1	A	407	VAL	CA-CB-CG1	6.07	120.00	110.90
1	A	467	SER	N-CA-C	6.03	127.27	111.00
1	A	52	ARG	CA-CB-CG	-6.03	100.14	113.40
1	A	326	TYR	CB-CG-CD1	5.96	124.58	121.00
1	A	408	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	276	ASP	CB-CG-OD1	5.84	123.55	118.30
1	A	392	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	435	LEU	CA-CB-CG	-5.75	102.08	115.30
1	A	115	ASP	CB-CG-OD2	5.74	123.47	118.30
1	A	372	ALA	C-N-CA	-5.70	107.45	121.70
1	A	85	ALA	N-CA-CB	5.67	118.04	110.10
1	A	249	ASP	O-C-N	5.65	131.73	122.70
1	A	473	GLY	N-CA-C	-5.63	99.02	113.10
1	A	177	ASP	CB-CG-OD2	5.63	123.36	118.30
1	A	401	ASP	CB-CG-OD1	-5.61	113.25	118.30
1	A	373	ASP	N-CA-CB	5.58	120.65	110.60
1	A	401	ASP	CB-CG-OD2	5.48	123.23	118.30
1	A	45	ASP	CB-CG-OD1	5.47	123.22	118.30
1	A	460	ILE	CA-CB-CG1	5.45	121.36	111.00
1	A	326	TYR	CB-CG-CD2	-5.45	117.73	121.00
1	A	266	ASP	CB-CG-OD1	5.43	123.19	118.30
1	A	426	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	A	460	ILE	N-CA-CB	-5.33	98.53	110.80
1	A	169	LEU	CB-CG-CD1	-5.32	101.95	111.00
1	A	71	ALA	CB-CA-C	5.32	118.08	110.10
1	A	498	ASP	N-CA-CB	-5.32	101.03	110.60
1	A	30	GLN	CB-CA-C	-5.31	99.78	110.40
1	A	181	ASP	CB-CG-OD1	-5.29	113.54	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	130	LEU	C-N-CA	-5.27	108.53	121.70
1	A	22	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	47	VAL	CA-CB-CG1	-5.24	103.04	110.90
1	A	253	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	271	CYS	N-CA-CB	5.20	119.97	110.60
1	A	260	ASP	CB-CG-OD1	5.17	122.95	118.30
1	A	234	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	A	182	ARG	NE-CZ-NH2	5.11	122.86	120.30
1	A	385	LEU	CB-CA-C	5.09	119.88	110.20
1	A	373	ASP	C-N-CA	5.05	134.34	121.70
1	A	420	ARG	NE-CZ-NH2	-5.03	117.78	120.30
1	A	26	TYR	CB-CG-CD2	5.03	124.02	121.00
1	A	256	GLN	CB-CA-C	-5.02	100.36	110.40
1	A	497	TYR	CB-CG-CD1	-5.02	117.99	121.00
1	A	498	ASP	CB-CG-OD2	5.01	122.81	118.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	251	THR	CA
1	A	254	LEU	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	467	SER	Mainchain

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3800	0	3787	471	0
2	A	15	0	7	2	0
3	A	33	0	0	2	0
All	All	3848	0	3794	471	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 62.

All (471) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:LEU:HB3	1:A:160:PRO:HB2	1.32	1.12
1:A:335:GLU:O	1:A:336:GLN:HB2	1.41	1.08
1:A:82:VAL:HG23	1:A:151:PHE:HA	1.30	1.06
1:A:332:GLU:HG3	1:A:337:ARG:HD3	1.10	1.06
1:A:214:ALA:HB3	1:A:217:SER:HB2	1.39	1.04
1:A:332:GLU:CG	1:A:337:ARG:HD3	1.87	1.04
1:A:274:MET:HE2	1:A:287:PRO:HA	1.40	1.00
1:A:53:GLU:HG2	1:A:61:PHE:HB3	1.39	1.00
1:A:271:CYS:HB3	1:A:327:VAL:HG13	1.46	0.97
1:A:5:GLN:HG2	1:A:6:PRO:HD3	1.45	0.96
1:A:462:LEU:HB3	1:A:478:ALA:HB3	1.44	0.96
1:A:429:HIS:HB3	1:A:430:PRO:HD2	1.49	0.93
1:A:32:THR:HG22	1:A:51:LYS:HE3	1.50	0.91
1:A:157:PHE:HE2	1:A:248:GLY:HA2	1.33	0.91
1:A:101:LEU:HB2	1:A:103:VAL:HG22	1.49	0.91
1:A:64:ARG:NH1	1:A:168:THR:HG22	1.86	0.90
1:A:340:LEU:H	1:A:409:LEU:HB3	1.38	0.89
1:A:435:LEU:H	1:A:503:THR:HG22	1.35	0.89
1:A:247:ILE:HD12	1:A:248:GLY:C	1.93	0.88
1:A:392:ARG:HH22	1:A:410:SER:HB2	1.38	0.88
1:A:347:GLU:HG3	1:A:350:GLY:H	1.38	0.88
1:A:358:LEU:HD12	1:A:358:LEU:H	1.39	0.87
1:A:79:ALA:HB1	1:A:103:VAL:CG1	2.04	0.86
1:A:79:ALA:HB1	1:A:103:VAL:HG12	1.55	0.86
1:A:174:LEU:HD13	1:A:204:MET:HE2	1.59	0.84
1:A:347:GLU:HG3	1:A:348:GLU:H	1.41	0.84
1:A:340:LEU:HB2	1:A:409:LEU:HB3	1.61	0.83
1:A:332:GLU:HG3	1:A:337:ARG:CD	2.03	0.82
1:A:149:GLN:HB2	1:A:151:PHE:CE1	2.16	0.81
1:A:159:HIS:CD2	1:A:160:PRO:HD2	2.16	0.80
1:A:158:ASP:HA	1:A:252:PHE:CE1	2.15	0.80
1:A:32:THR:CG2	1:A:51:LYS:HE3	2.12	0.80
1:A:174:LEU:HD21	1:A:200:ILE:HD13	1.62	0.80
1:A:245:LYS:HG2	1:A:246:ARG:N	1.94	0.80
1:A:340:LEU:HB2	1:A:409:LEU:CB	2.11	0.80
1:A:508:PHE:CD1	1:A:512:LEU:HB2	2.18	0.79
1:A:462:LEU:HB3	1:A:478:ALA:CB	2.12	0.79
1:A:7:LEU:HD12	1:A:161:MET:SD	2.23	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:VAL:HG21	1:A:211:ALA:HB1	1.64	0.79
1:A:159:HIS:HD2	1:A:161:MET:H	1.31	0.78
1:A:337:ARG:HH22	1:A:416:LYS:HD2	1.48	0.78
1:A:392:ARG:NH2	1:A:410:SER:HB2	1.97	0.78
1:A:308:GLU:HB2	1:A:310:LEU:HD21	1.66	0.77
1:A:247:ILE:HD12	1:A:248:GLY:N	2.00	0.77
1:A:101:LEU:HB2	1:A:103:VAL:CG2	2.15	0.77
1:A:7:LEU:HB3	1:A:160:PRO:CB	2.12	0.76
1:A:20:VAL:HG11	1:A:199:LEU:HD21	1.67	0.76
1:A:228:HIS:O	1:A:230:VAL:HG23	1.84	0.76
1:A:340:LEU:N	1:A:409:LEU:HB3	2.01	0.76
1:A:219:CYS:HB2	3:A:601:HOH:O	1.88	0.74
1:A:394:GLU:O	1:A:395:ILE:C	2.26	0.74
1:A:435:LEU:H	1:A:503:THR:CG2	2.00	0.74
1:A:6:PRO:O	1:A:7:LEU:HG	1.87	0.74
1:A:366:GLU:HA	1:A:465:TYR:O	1.88	0.74
1:A:289:GLY:HA2	1:A:315:SER:OG	1.87	0.73
1:A:409:LEU:HD13	1:A:410:SER:N	2.02	0.73
1:A:447:LEU:O	1:A:448:LEU:C	2.23	0.73
1:A:460:ILE:HD12	1:A:479:PHE:HE1	1.51	0.73
1:A:438:PHE:HA	1:A:498:ASP:O	1.89	0.72
1:A:194:ALA:O	1:A:198:VAL:HG23	1.89	0.71
1:A:271:CYS:HB3	1:A:327:VAL:CG1	2.20	0.71
1:A:344:THR:HG22	1:A:376:ASN:HB3	1.71	0.71
1:A:165:GLY:O	1:A:168:THR:HB	1.91	0.71
1:A:187:VAL:CG2	1:A:211:ALA:HB1	2.21	0.71
1:A:248:GLY:HA3	1:A:252:PHE:CG	2.26	0.70
1:A:347:GLU:CG	1:A:350:GLY:H	2.05	0.70
1:A:90:HIS:CE1	1:A:156:PRO:HA	2.26	0.70
1:A:274:MET:HE1	1:A:287:PRO:HG3	1.72	0.70
1:A:17:LEU:HD11	1:A:203:LEU:CD1	2.22	0.70
1:A:157:PHE:CE2	1:A:248:GLY:HA2	2.22	0.69
1:A:375:LYS:HD2	1:A:375:LYS:O	1.92	0.69
1:A:229:PRO:HB2	1:A:246:ARG:CZ	2.22	0.69
1:A:7:LEU:CD2	1:A:11:PRO:HG3	2.23	0.69
1:A:46:ASN:HB3	1:A:308:GLU:O	1.93	0.69
1:A:462:LEU:CB	1:A:478:ALA:HB3	2.22	0.69
1:A:247:ILE:HG23	3:A:616:HOH:O	1.91	0.69
1:A:358:LEU:HD12	1:A:358:LEU:N	2.07	0.69
1:A:62:LYS:HD2	1:A:90:HIS:HB2	1.76	0.68
1:A:215:GLU:OE1	1:A:266:ASP:HB3	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:GLU:O	1:A:434:ARG:HG2	1.94	0.68
1:A:186:PRO:HD3	1:A:292:ALA:HB2	1.76	0.67
1:A:365:THR:O	1:A:366:GLU:HG2	1.94	0.67
1:A:337:ARG:NH2	1:A:416:LYS:HD2	2.10	0.67
1:A:394:GLU:O	1:A:396:LEU:N	2.28	0.67
1:A:472:TYR:HA	1:A:474:ARG:HH12	1.56	0.67
1:A:274:MET:HE2	1:A:287:PRO:CA	2.21	0.67
1:A:335:GLU:O	1:A:336:GLN:CB	2.24	0.67
1:A:392:ARG:NE	1:A:409:LEU:HD11	2.10	0.67
1:A:500:HIS:ND1	1:A:502:GLU:OE2	2.28	0.67
1:A:291:LEU:HD12	1:A:291:LEU:O	1.95	0.66
1:A:247:ILE:HD12	1:A:248:GLY:CA	2.24	0.66
1:A:345:ILE:HD12	1:A:379:ILE:HG13	1.77	0.66
1:A:7:LEU:CB	1:A:160:PRO:HB2	2.18	0.66
1:A:7:LEU:HD22	1:A:11:PRO:HG3	1.76	0.66
1:A:141:LYS:NZ	1:A:144:GLU:OE2	2.28	0.66
1:A:135:PHE:O	1:A:138:ALA:N	2.29	0.66
1:A:182:ARG:NH1	1:A:308:GLU:OE2	2.29	0.66
1:A:158:ASP:OD2	1:A:249:ASP:HB2	1.96	0.65
1:A:472:TYR:HA	1:A:474:ARG:NH1	2.11	0.65
1:A:286:GLU:HB2	1:A:319:VAL:HG22	1.79	0.65
1:A:143:ILE:O	1:A:146:SER:HB2	1.97	0.65
1:A:228:HIS:O	1:A:229:PRO:C	2.33	0.65
1:A:395:ILE:HG22	1:A:396:LEU:N	2.12	0.65
1:A:425:GLY:O	1:A:512:LEU:HA	1.95	0.65
1:A:434:ARG:NH2	1:A:501:ASP:OD2	2.30	0.65
1:A:250:GLU:O	1:A:254:LEU:N	2.30	0.64
1:A:396:LEU:HD11	1:A:407:VAL:CG2	2.27	0.64
1:A:219:CYS:HB3	1:A:247:ILE:HG22	1.78	0.64
1:A:262:ILE:HG13	1:A:263:ILE:N	2.11	0.64
1:A:426:ARG:NH2	1:A:511:PHE:HA	2.13	0.64
1:A:11:PRO:HB2	1:A:16:TYR:CE2	2.33	0.64
1:A:321:PHE:O	1:A:324:LEU:HB2	1.97	0.64
1:A:78:LYS:O	1:A:78:LYS:HG2	1.98	0.63
1:A:270:ILE:HG22	1:A:274:MET:CE	2.29	0.63
1:A:449:ARG:O	1:A:450:PHE:C	2.36	0.63
1:A:257:GLU:OE1	1:A:257:GLU:HA	1.98	0.63
1:A:440:PHE:H	1:A:474:ARG:HG3	1.64	0.63
1:A:353:LEU:O	1:A:356:CYS:N	2.32	0.63
1:A:392:ARG:NH1	1:A:409:LEU:HD21	2.14	0.63
1:A:447:LEU:HD22	1:A:451:LEU:HD22	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:GLN:HG2	1:A:6:PRO:CD	2.25	0.62
1:A:64:ARG:HH11	1:A:168:THR:HG22	1.64	0.62
1:A:174:LEU:HD13	1:A:204:MET:CE	2.27	0.62
1:A:340:LEU:H	1:A:409:LEU:CB	2.10	0.62
1:A:435:LEU:N	1:A:503:THR:HG22	2.11	0.62
1:A:215:GLU:HB2	1:A:265:VAL:O	1.99	0.62
1:A:121:VAL:O	1:A:122:ARG:C	2.38	0.61
1:A:346:PRO:HB3	1:A:376:ASN:OD1	2.01	0.61
1:A:293:LEU:HD12	1:A:293:LEU:O	2.01	0.61
1:A:378:CYS:C	1:A:379:ILE:HG12	2.19	0.61
1:A:392:ARG:CZ	1:A:409:LEU:HD11	2.30	0.61
1:A:286:GLU:O	1:A:287:PRO:C	2.37	0.61
1:A:252:PHE:O	1:A:253:ARG:C	2.36	0.60
1:A:387:ARG:NH2	1:A:394:GLU:OE1	2.33	0.60
1:A:461:SER:N	1:A:478:ALA:O	2.30	0.60
1:A:82:VAL:HG21	1:A:151:PHE:CD2	2.36	0.60
1:A:286:GLU:CB	1:A:319:VAL:HG22	2.31	0.60
1:A:347:GLU:CG	1:A:348:GLU:H	2.11	0.60
1:A:347:GLU:HG2	1:A:350:GLY:HA3	1.84	0.60
1:A:508:PHE:CE2	1:A:513:ALA:HB2	2.37	0.60
1:A:336:GLN:HE22	1:A:389:LEU:HB2	1.66	0.60
1:A:20:VAL:HG21	1:A:199:LEU:HD22	1.83	0.60
1:A:107:ILE:HD13	1:A:121:VAL:CG1	2.32	0.60
1:A:434:ARG:HA	1:A:503:THR:HG22	1.82	0.60
1:A:442:GLU:OE1	1:A:467:SER:OG	2.19	0.59
1:A:418:HIS:O	1:A:421:TYR:HB2	2.03	0.59
1:A:64:ARG:HH12	1:A:169:LEU:HA	1.67	0.59
1:A:340:LEU:CB	1:A:409:LEU:HB3	2.32	0.59
1:A:157:PHE:O	1:A:252:PHE:HZ	1.85	0.59
1:A:395:ILE:HG22	1:A:396:LEU:HD22	1.84	0.59
1:A:98:SER:HB3	1:A:105:ALA:HB2	1.84	0.59
1:A:392:ARG:HH22	1:A:410:SER:CB	2.14	0.59
1:A:133:ALA:HB3	1:A:137:GLU:OE1	2.03	0.58
1:A:433:GLU:C	1:A:434:ARG:HG2	2.23	0.58
1:A:114:ALA:O	1:A:115:ASP:C	2.40	0.58
1:A:159:HIS:C	1:A:163:ILE:HD12	2.23	0.58
1:A:270:ILE:HG22	1:A:274:MET:HE2	1.85	0.58
1:A:17:LEU:HD11	1:A:203:LEU:HD11	1.85	0.58
1:A:17:LEU:O	1:A:20:VAL:HB	2.04	0.58
1:A:338:GLU:OE2	1:A:420:ARG:NH2	2.29	0.58
1:A:52:ARG:HG2	1:A:54:ASP:OD1	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:MET:O	1:A:297:LYS:C	2.41	0.58
1:A:339:ALA:HA	1:A:409:LEU:HG	1.86	0.58
1:A:212:VAL:HG22	1:A:263:ILE:HG22	1.85	0.57
1:A:317:ALA:O	1:A:319:VAL:HG23	2.04	0.57
1:A:20:VAL:HG11	1:A:199:LEU:CD2	2.33	0.57
1:A:199:LEU:O	1:A:200:ILE:C	2.40	0.57
1:A:407:VAL:HG13	1:A:409:LEU:CD1	2.34	0.57
1:A:252:PHE:CD1	1:A:252:PHE:N	2.70	0.57
1:A:81:GLY:C	1:A:82:VAL:HG22	2.25	0.57
1:A:447:LEU:O	1:A:450:PHE:N	2.37	0.57
1:A:508:PHE:O	1:A:513:ALA:N	2.29	0.57
1:A:53:GLU:OE2	1:A:56:GLN:NE2	2.36	0.57
1:A:86:SER:HA	1:A:108:VAL:O	2.05	0.57
1:A:141:LYS:O	1:A:144:GLU:HB3	2.05	0.56
1:A:158:ASP:HA	1:A:252:PHE:CZ	2.39	0.56
1:A:214:ALA:O	1:A:215:GLU:C	2.42	0.56
1:A:181:ASP:OD2	1:A:308:GLU:HB3	2.05	0.56
1:A:223:ALA:O	1:A:227:GLY:N	2.37	0.56
1:A:7:LEU:O	1:A:9:GLY:N	2.29	0.56
1:A:157:PHE:HD2	1:A:252:PHE:CE2	2.23	0.56
1:A:462:LEU:HD23	1:A:463:PHE:N	2.20	0.56
1:A:146:SER:O	1:A:147:GLN:C	2.40	0.56
1:A:10:ALA:O	1:A:11:PRO:C	2.44	0.56
1:A:116:ILE:HG23	1:A:117:LYS:N	2.20	0.56
1:A:220:LEU:O	1:A:221:LYS:C	2.44	0.56
1:A:157:PHE:HE2	1:A:248:GLY:CA	2.12	0.56
1:A:162:VAL:O	1:A:166:GLN:HG2	2.05	0.56
1:A:339:ALA:O	1:A:382:GLY:HA2	2.06	0.56
1:A:332:GLU:CD	1:A:337:ARG:HD3	2.25	0.56
1:A:53:GLU:HG2	1:A:61:PHE:CB	2.25	0.55
1:A:286:GLU:O	1:A:289:GLY:N	2.39	0.55
1:A:460:ILE:HD12	1:A:479:PHE:CE1	2.37	0.55
1:A:47:VAL:O	1:A:310:LEU:N	2.28	0.55
1:A:286:GLU:OE1	1:A:318:ASN:N	2.39	0.55
1:A:70:MET:HE2	1:A:101:LEU:HD12	1.87	0.55
1:A:81:GLY:O	1:A:82:VAL:HG13	2.06	0.55
1:A:179:HIS:O	1:A:309:ARG:HD3	2.06	0.55
1:A:187:VAL:HG21	1:A:211:ALA:CB	2.36	0.55
1:A:247:ILE:O	1:A:248:GLY:O	2.25	0.55
1:A:509:ARG:HA	1:A:513:ALA:HB3	1.89	0.55
1:A:107:ILE:CG2	1:A:128:VAL:HG12	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:VAL:CG1	1:A:243:ALA:HB1	2.36	0.55
1:A:14:ALA:O	1:A:15:GLU:C	2.44	0.55
1:A:249:ASP:O	1:A:253:ARG:N	2.39	0.55
1:A:442:GLU:HG2	1:A:468:HIS:ND1	2.22	0.55
1:A:76:GLU:HB3	1:A:78:LYS:HE2	1.87	0.54
1:A:397:GLN:O	1:A:400:ASN:N	2.37	0.54
1:A:36:LYS:HE3	1:A:38:GLU:OE2	2.06	0.54
1:A:107:ILE:HD13	1:A:121:VAL:HG11	1.88	0.54
1:A:301:ALA:O	1:A:302:LEU:C	2.43	0.54
1:A:61:PHE:O	1:A:64:ARG:HG3	2.08	0.54
1:A:70:MET:CE	1:A:83:ILE:HD13	2.38	0.54
1:A:213:GLU:O	1:A:264:THR:HA	2.07	0.54
1:A:74:THR:CG2	1:A:79:ALA:HB2	2.38	0.54
1:A:197:ALA:O	1:A:198:VAL:C	2.46	0.54
1:A:284:VAL:O	1:A:284:VAL:HG12	2.08	0.54
1:A:391:GLU:HA	1:A:394:GLU:HG3	1.90	0.54
1:A:462:LEU:HD23	1:A:462:LEU:C	2.28	0.54
1:A:291:LEU:HD12	1:A:291:LEU:C	2.28	0.53
1:A:296:MET:O	1:A:299:TYR:N	2.41	0.53
1:A:56:GLN:HG3	1:A:60:SER:O	2.08	0.53
1:A:426:ARG:NH2	1:A:510:PHE:O	2.36	0.53
1:A:233:PRO:HG2	1:A:234:ARG:H	1.73	0.53
1:A:342:ALA:HB3	1:A:406:VAL:HB	1.91	0.53
1:A:246:ARG:C	1:A:247:ILE:O	2.44	0.52
1:A:475:VAL:HG13	1:A:476:LEU:N	2.24	0.52
1:A:11:PRO:HG2	1:A:16:TYR:CZ	2.45	0.52
1:A:15:GLU:O	1:A:18:ARG:HB3	2.08	0.52
1:A:347:GLU:OE1	1:A:347:GLU:HA	2.09	0.52
1:A:17:LEU:HD23	1:A:17:LEU:C	2.30	0.52
1:A:392:ARG:O	1:A:396:LEU:HD23	2.10	0.52
1:A:439:GLU:OE1	1:A:472:TYR:O	2.28	0.52
1:A:81:GLY:O	1:A:82:VAL:HG22	2.10	0.52
1:A:107:ILE:HB	1:A:128:VAL:HG12	1.90	0.52
1:A:70:MET:HE3	1:A:83:ILE:HD13	1.91	0.52
1:A:185:VAL:HG21	1:A:193:ALA:HA	1.92	0.52
1:A:309:ARG:CG	1:A:309:ARG:HH11	2.23	0.52
1:A:16:TYR:OH	1:A:160:PRO:HB3	2.09	0.51
1:A:245:LYS:HG2	1:A:246:ARG:H	1.72	0.51
1:A:246:ARG:O	1:A:247:ILE:O	2.28	0.51
1:A:472:TYR:CE1	1:A:474:ARG:NH2	2.78	0.51
1:A:111:THR:HG23	1:A:132:GLY:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:508:PHE:O	1:A:512:LEU:N	2.38	0.51
1:A:157:PHE:O	1:A:159:HIS:N	2.44	0.51
1:A:37:MET:O	1:A:41:SER:HB2	2.10	0.51
1:A:41:SER:HA	1:A:48:ILE:HG13	1.93	0.51
1:A:53:GLU:CD	1:A:56:GLN:HE21	2.13	0.51
1:A:14:ALA:O	1:A:17:LEU:N	2.44	0.51
1:A:107:ILE:HG22	1:A:107:ILE:O	2.11	0.51
1:A:7:LEU:HD22	1:A:11:PRO:CG	2.40	0.51
1:A:89:ASN:ND2	2:A:962:PLP:O3	2.45	0.50
1:A:396:LEU:HD11	1:A:407:VAL:HG22	1.92	0.50
1:A:51:LYS:O	1:A:53:GLU:N	2.42	0.50
1:A:122:ARG:O	1:A:125:GLY:O	2.29	0.50
1:A:295:GLY:O	1:A:296:MET:C	2.49	0.50
1:A:295:GLY:O	1:A:298:LYS:N	2.44	0.50
1:A:352:PHE:N	1:A:352:PHE:CD1	2.79	0.50
1:A:370:ARG:O	1:A:371:PHE:C	2.50	0.50
1:A:250:GLU:O	1:A:252:PHE:N	2.44	0.50
1:A:302:LEU:HB3	1:A:303:HIS:ND1	2.26	0.50
1:A:340:LEU:CA	1:A:409:LEU:HB3	2.42	0.50
1:A:12:GLU:O	1:A:13:GLY:C	2.50	0.50
1:A:212:VAL:HG12	1:A:291:LEU:HD11	1.93	0.50
1:A:167:GLY:O	1:A:170:ALA:N	2.45	0.49
1:A:82:VAL:CG2	1:A:151:PHE:CD2	2.95	0.49
1:A:449:ARG:HB3	1:A:497:TYR:OH	2.13	0.49
1:A:269:ALA:O	1:A:270:ILE:C	2.51	0.49
1:A:44:LEU:HD13	1:A:300:ILE:HD12	1.94	0.49
1:A:82:VAL:HG23	1:A:151:PHE:CA	2.23	0.49
1:A:141:LYS:O	1:A:142:ALA:C	2.51	0.49
1:A:255:CYS:C	1:A:257:GLU:H	2.15	0.49
1:A:266:ASP:O	1:A:269:ALA:N	2.45	0.49
1:A:35:GLN:OE1	1:A:52:ARG:NH1	2.46	0.49
1:A:145:LEU:O	1:A:146:SER:O	2.31	0.49
1:A:157:PHE:CZ	1:A:189:GLY:HA3	2.48	0.49
1:A:417:LEU:O	1:A:417:LEU:HG	2.10	0.49
1:A:139:LYS:HG3	1:A:143:ILE:HD12	1.95	0.48
1:A:84:THR:HG23	1:A:84:THR:O	2.11	0.48
1:A:433:GLU:HA	1:A:479:PHE:O	2.12	0.48
1:A:182:ARG:HG3	1:A:310:LEU:CD2	2.44	0.48
1:A:256:GLN:N	1:A:256:GLN:OE1	2.47	0.48
1:A:131:HIS:O	1:A:138:ALA:HA	2.13	0.48
1:A:437:SER:OG	1:A:502:GLU:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:TYR:HE1	1:A:458:TRP:CZ2	2.32	0.48
1:A:252:PHE:O	1:A:256:GLN:OE1	2.32	0.48
1:A:157:PHE:CE1	1:A:189:GLY:HA3	2.48	0.48
1:A:451:LEU:N	1:A:451:LEU:CD1	2.77	0.48
1:A:5:GLN:O	1:A:7:LEU:N	2.47	0.47
1:A:199:LEU:HD13	1:A:199:LEU:HA	1.64	0.47
1:A:396:LEU:N	1:A:396:LEU:HD22	2.29	0.47
1:A:508:PHE:CD2	1:A:513:ALA:HB2	2.49	0.47
1:A:63:LEU:HD12	1:A:63:LEU:O	2.14	0.47
1:A:75:GLU:O	1:A:75:GLU:HG2	2.13	0.47
1:A:374:ALA:O	1:A:375:LYS:C	2.53	0.47
1:A:409:LEU:HD13	1:A:409:LEU:C	2.34	0.47
1:A:448:LEU:O	1:A:449:ARG:C	2.51	0.47
1:A:191:GLY:N	1:A:252:PHE:HE2	2.12	0.47
1:A:434:ARG:HE	1:A:434:ARG:HB3	1.47	0.47
1:A:7:LEU:HD23	1:A:11:PRO:HG3	1.95	0.47
1:A:232:LEU:HD22	1:A:233:PRO:HD2	1.95	0.47
1:A:261:ASP:OD1	1:A:262:ILE:N	2.37	0.47
1:A:426:ARG:NH1	1:A:513:ALA:O	2.48	0.47
1:A:435:LEU:HD12	1:A:477:ALA:O	2.13	0.47
1:A:457:TYR:CE1	1:A:458:TRP:CZ2	3.02	0.47
1:A:347:GLU:HG3	1:A:350:GLY:N	2.19	0.47
1:A:186:PRO:HD3	1:A:292:ALA:CB	2.43	0.47
1:A:293:LEU:HD12	1:A:293:LEU:C	2.34	0.47
1:A:385:LEU:HD11	1:A:392:ARG:HB2	1.96	0.47
1:A:431:LEU:HD12	1:A:431:LEU:HA	1.64	0.47
1:A:64:ARG:NH1	1:A:169:LEU:CA	2.77	0.47
1:A:84:THR:HA	1:A:106:LEU:O	2.14	0.47
1:A:182:ARG:HG3	1:A:310:LEU:HD22	1.96	0.47
1:A:270:ILE:HG22	1:A:274:MET:HE3	1.96	0.47
1:A:385:LEU:HB2	1:A:391:GLU:OE1	2.15	0.47
1:A:74:THR:HG23	1:A:79:ALA:HB2	1.96	0.47
1:A:145:LEU:HA	1:A:145:LEU:HD23	1.55	0.47
1:A:451:LEU:N	1:A:451:LEU:HD13	2.30	0.47
1:A:86:SER:O	1:A:87:ALA:HB2	2.14	0.47
1:A:18:ARG:O	1:A:22:ARG:HB2	2.15	0.46
1:A:200:ILE:CG2	1:A:207:ILE:HB	2.46	0.46
1:A:213:GLU:OE2	1:A:220:LEU:HB3	2.15	0.46
1:A:417:LEU:CD2	1:A:418:HIS:CD2	2.98	0.46
1:A:381:VAL:CG1	1:A:382:GLY:N	2.79	0.46
1:A:69:MET:CE	1:A:154:VAL:HG22	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:SER:HA	1:A:103:VAL:HG23	1.96	0.46
1:A:187:VAL:HG11	1:A:220:LEU:HD22	1.97	0.46
1:A:309:ARG:NH1	1:A:309:ARG:HG2	2.29	0.46
1:A:497:TYR:CD1	1:A:497:TYR:N	2.83	0.46
1:A:171:LEU:HD23	1:A:171:LEU:HA	1.55	0.46
1:A:392:ARG:HH11	1:A:409:LEU:HD21	1.79	0.46
1:A:447:LEU:HA	1:A:447:LEU:HD23	1.70	0.45
1:A:12:GLU:O	1:A:15:GLU:N	2.50	0.45
1:A:46:ASN:OD1	1:A:305:ILE:HG22	2.16	0.45
1:A:76:GLU:OE1	1:A:78:LYS:HD3	2.16	0.45
1:A:229:PRO:O	1:A:246:ARG:NH1	2.49	0.45
1:A:407:VAL:CG1	1:A:409:LEU:CD1	2.94	0.45
1:A:10:ALA:N	1:A:11:PRO:CD	2.79	0.45
1:A:385:LEU:HD13	1:A:391:GLU:HG2	1.98	0.45
1:A:447:LEU:O	1:A:450:PHE:HB3	2.17	0.45
1:A:116:ILE:CG2	1:A:117:LYS:N	2.79	0.45
1:A:185:VAL:HA	1:A:186:PRO:HD2	1.89	0.45
1:A:429:HIS:HB3	1:A:430:PRO:CD	2.34	0.45
1:A:440:PHE:O	1:A:441:PRO:C	2.55	0.45
1:A:11:PRO:HG2	1:A:16:TYR:OH	2.16	0.45
1:A:107:ILE:CB	1:A:128:VAL:HG12	2.47	0.45
1:A:212:VAL:HA	1:A:263:ILE:O	2.17	0.45
1:A:407:VAL:HG13	1:A:409:LEU:HD12	1.98	0.45
1:A:415:ALA:HA	1:A:419:VAL:HB	1.98	0.45
1:A:69:MET:HE2	1:A:154:VAL:HG22	1.97	0.45
1:A:249:ASP:H	1:A:252:PHE:HD1	1.63	0.45
1:A:16:TYR:O	1:A:17:LEU:C	2.54	0.45
1:A:157:PHE:CD2	1:A:252:PHE:CE2	3.05	0.45
1:A:356:CYS:O	1:A:359:LEU:N	2.44	0.45
1:A:448:LEU:HD12	1:A:448:LEU:HA	1.62	0.45
1:A:391:GLU:O	1:A:394:GLU:HG3	2.16	0.44
1:A:107:ILE:HD13	1:A:121:VAL:HG13	1.98	0.44
1:A:296:MET:HB2	1:A:296:MET:HE3	1.71	0.44
1:A:351:SER:OG	1:A:352:PHE:N	2.47	0.44
1:A:387:ARG:O	1:A:388:GLY:C	2.55	0.44
1:A:159:HIS:HB3	1:A:162:VAL:HB	2.00	0.44
1:A:324:LEU:HD13	1:A:324:LEU:HA	1.47	0.44
1:A:409:LEU:HD22	1:A:409:LEU:O	2.17	0.44
1:A:5:GLN:N	1:A:6:PRO:CD	2.80	0.44
1:A:66:ALA:O	1:A:70:MET:HG3	2.17	0.44
1:A:154:VAL:HA	1:A:155:PRO:HD2	1.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:PHE:HB3	1:A:239:ALA:H	1.60	0.44
1:A:352:PHE:CE2	1:A:379:ILE:CD1	3.00	0.44
1:A:409:LEU:HD12	1:A:409:LEU:H	1.83	0.44
1:A:409:LEU:CD1	1:A:409:LEU:H	2.31	0.44
1:A:141:LYS:HD2	1:A:141:LYS:HA	1.64	0.44
1:A:5:GLN:N	1:A:6:PRO:HD2	2.32	0.44
1:A:233:PRO:CG	1:A:234:ARG:H	2.30	0.44
1:A:358:LEU:N	1:A:358:LEU:CD1	2.80	0.44
1:A:98:SER:HB2	1:A:103:VAL:O	2.18	0.44
1:A:293:LEU:O	1:A:296:MET:HB3	2.18	0.44
1:A:429:HIS:O	1:A:430:PRO:C	2.56	0.44
1:A:251:THR:O	1:A:255:CYS:HB2	2.17	0.43
1:A:512:LEU:O	1:A:513:ALA:C	2.52	0.43
1:A:190:GLY:HA3	1:A:252:PHE:HD2	1.83	0.43
1:A:351:SER:OG	1:A:353:LEU:HD23	2.17	0.43
1:A:380:PHE:CG	1:A:423:VAL:HG21	2.53	0.43
1:A:407:VAL:HG13	1:A:409:LEU:HD11	1.99	0.43
1:A:119:ASP:O	1:A:123:GLY:N	2.46	0.43
1:A:83:ILE:HA	1:A:152:THR:H	1.84	0.43
1:A:189:GLY:HA2	1:A:244:VAL:HG11	1.99	0.43
1:A:409:LEU:C	1:A:409:LEU:HD22	2.38	0.43
1:A:434:ARG:HB2	1:A:436:TYR:HE1	1.82	0.43
1:A:268:ASP:OD2	1:A:330:ARG:NH1	2.51	0.43
1:A:398:MET:O	1:A:400:ASN:N	2.52	0.43
1:A:419:VAL:H	1:A:419:VAL:HG23	1.57	0.43
1:A:448:LEU:O	1:A:451:LEU:N	2.51	0.43
1:A:247:ILE:CD1	1:A:248:GLY:N	2.76	0.43
1:A:434:ARG:NE	1:A:503:THR:HG21	2.34	0.43
1:A:443:SER:OG	1:A:444:PRO:HD2	2.19	0.43
1:A:79:ALA:HB1	1:A:103:VAL:HG11	1.94	0.43
1:A:390:GLU:H	1:A:390:GLU:HG3	1.64	0.43
1:A:250:GLU:O	1:A:251:THR:C	2.57	0.43
1:A:331:CYS:O	1:A:332:GLU:C	2.57	0.43
1:A:352:PHE:O	1:A:355:PHE:HB3	2.18	0.43
1:A:158:ASP:OD2	1:A:249:ASP:N	2.51	0.43
1:A:358:LEU:H	1:A:358:LEU:CD1	2.20	0.43
1:A:396:LEU:HD13	1:A:396:LEU:HA	1.42	0.43
1:A:503:THR:C	1:A:505:ASN:H	2.21	0.43
1:A:64:ARG:NH1	1:A:169:LEU:HA	2.34	0.43
1:A:74:THR:CG2	1:A:79:ALA:CB	2.97	0.43
1:A:76:GLU:O	1:A:78:LYS:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HA	1:A:247:ILE:HG21	2.00	0.43
1:A:173:LEU:HD12	1:A:173:LEU:HA	1.66	0.42
1:A:347:GLU:HG2	1:A:350:GLY:CA	2.49	0.42
1:A:435:LEU:HG	1:A:476:LEU:HD21	2.01	0.42
1:A:108:VAL:CG2	1:A:142:ALA:HB2	2.49	0.42
1:A:381:VAL:HG12	1:A:382:GLY:N	2.34	0.42
1:A:64:ARG:HH12	1:A:169:LEU:CA	2.32	0.42
1:A:198:VAL:O	1:A:202:GLN:HG2	2.19	0.42
1:A:229:PRO:HB2	1:A:246:ARG:NE	2.35	0.42
1:A:47:VAL:N	1:A:308:GLU:O	2.48	0.42
1:A:74:THR:HG22	1:A:79:ALA:CB	2.50	0.42
1:A:343:VAL:HG13	1:A:343:VAL:O	2.20	0.42
1:A:52:ARG:HH11	1:A:52:ARG:HD3	1.34	0.42
1:A:199:LEU:O	1:A:202:GLN:N	2.50	0.42
1:A:400:ASN:O	1:A:400:ASN:ND2	2.53	0.42
1:A:70:MET:CE	1:A:83:ILE:CD1	2.97	0.42
1:A:272:ALA:O	1:A:275:LYS:N	2.52	0.42
1:A:344:THR:HG22	1:A:376:ASN:CB	2.47	0.42
1:A:344:THR:HA	1:A:378:CYS:HA	2.01	0.42
1:A:459:ASN:CB	1:A:480:GLU:HB2	2.50	0.42
1:A:142:ALA:CB	1:A:153:TRP:HZ3	2.32	0.42
1:A:204:MET:HA	1:A:205:PRO:HD2	1.74	0.42
1:A:220:LEU:HG	1:A:224:LEU:HD22	2.02	0.42
1:A:271:CYS:CB	1:A:327:VAL:HG13	2.33	0.42
1:A:16:TYR:OH	1:A:160:PRO:HA	2.19	0.42
1:A:328:SER:O	1:A:329:GLU:C	2.59	0.42
1:A:442:GLU:HG3	1:A:443:SER:N	2.34	0.42
1:A:106:LEU:HG	1:A:107:ILE:N	2.34	0.42
1:A:347:GLU:CG	1:A:348:GLU:N	2.80	0.42
1:A:5:GLN:CG	1:A:6:PRO:HD3	2.33	0.41
1:A:201:LYS:HD2	1:A:201:LYS:HA	1.64	0.41
1:A:436:TYR:O	1:A:476:LEU:HD23	2.19	0.41
1:A:465:TYR:HD1	1:A:475:VAL:HG23	1.86	0.41
1:A:83:ILE:HG22	1:A:152:THR:HB	2.02	0.41
1:A:130:LEU:HA	1:A:130:LEU:HD12	1.64	0.41
1:A:157:PHE:CD1	1:A:189:GLY:HA3	2.56	0.41
1:A:220:LEU:O	1:A:222:ALA:N	2.53	0.41
1:A:172:GLU:O	1:A:173:LEU:C	2.58	0.41
1:A:330:ARG:HD2	1:A:330:ARG:HA	1.86	0.41
1:A:409:LEU:CD1	1:A:409:LEU:N	2.84	0.41
1:A:299:TYR:CE1	1:A:303:HIS:CE1	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:LEU:C	1:A:513:ALA:O	2.55	0.41
1:A:345:ILE:HA	1:A:346:PRO:HD3	1.91	0.41
1:A:21:LEU:HD23	1:A:21:LEU:HA	1.86	0.41
1:A:177:ASP:CG	1:A:309:ARG:HE	2.24	0.41
1:A:189:GLY:N	1:A:244:VAL:HG21	2.36	0.41
1:A:242:VAL:CG2	1:A:288:SER:HB3	2.51	0.41
1:A:250:GLU:C	1:A:252:PHE:N	2.72	0.41
1:A:277:LEU:HD12	1:A:277:LEU:HA	1.61	0.41
1:A:343:VAL:H	1:A:343:VAL:HG12	1.56	0.41
1:A:419:VAL:HA	1:A:422:MET:HB2	2.02	0.41
1:A:476:LEU:HD23	1:A:476:LEU:HA	1.84	0.41
1:A:17:LEU:O	1:A:17:LEU:HD23	2.21	0.41
1:A:36:LYS:HE3	1:A:38:GLU:CD	2.41	0.41
1:A:107:ILE:HG22	1:A:128:VAL:HG12	2.01	0.41
1:A:159:HIS:CD2	1:A:161:MET:H	2.21	0.41
1:A:220:LEU:O	1:A:223:ALA:N	2.54	0.40
1:A:417:LEU:HD21	1:A:418:HIS:CD2	2.57	0.40
1:A:417:LEU:HD23	1:A:418:HIS:CD2	2.57	0.40
1:A:508:PHE:CE1	1:A:512:LEU:HB2	2.56	0.40
1:A:116:ILE:HG23	1:A:117:LYS:HG2	2.03	0.40
1:A:344:THR:CG2	1:A:376:ASN:HB3	2.46	0.40
1:A:224:LEU:HD12	1:A:224:LEU:HA	1.83	0.40
1:A:316:GLY:N	2:A:962:PLP:H2A2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	490/514 (95%)	360 (74%)	91 (19%)	39 (8%)	1 2

All (39) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	PRO
1	A	8	SER
1	A	52	ARG
1	A	82	VAL
1	A	114	ALA
1	A	126	GLY
1	A	146	SER
1	A	247	ILE
1	A	248	GLY
1	A	395	ILE
1	A	13	GLY
1	A	62	LYS
1	A	133	ALA
1	A	158	ASP
1	A	236	GLY
1	A	336	GLN
1	A	357	GLN
1	A	427	PRO
1	A	7	LEU
1	A	88	GLY
1	A	145	LEU
1	A	199	LEU
1	A	347	GLU
1	A	348	GLU
1	A	396	LEU
1	A	509	ARG
1	A	59	HIS
1	A	144	GLU
1	A	147	GLN
1	A	81	GLY
1	A	87	ALA
1	A	115	ASP
1	A	256	GLN
1	A	353	LEU
1	A	394	GLU
1	A	430	PRO
1	A	11	PRO
1	A	287	PRO
1	A	110	PRO

5.3.2 Protein sidechains (i)

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	391/408 (96%)	294 (75%)	97 (25%)	0 2

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	8	SER
1	A	17	LEU
1	A	18	ARG
1	A	22	ARG
1	A	24	PRO
1	A	30	GLN
1	A	32	THR
1	A	34	LEU
1	A	36	LYS
1	A	43	ARG
1	A	46	ASN
1	A	49	LEU
1	A	52	ARG
1	A	58	VAL
1	A	69	MET
1	A	74	THR
1	A	77	GLN
1	A	78	LYS
1	A	82	VAL
1	A	86	SER
1	A	100	ARG
1	A	103	VAL
1	A	108	VAL
1	A	111	THR
1	A	115	ASP
1	A	134	ASN
1	A	146	SER
1	A	148	GLN
1	A	158	ASP

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Mol	Chain	Res	Type
1	A	168	THR
1	A	171	LEU
1	A	175	GLN
1	A	182	ARG
1	A	199	LEU
1	A	207	ILE
1	A	208	LYS
1	A	216	ASP
1	A	217	SER
1	A	219	CYS
1	A	224	LEU
1	A	228	HIS
1	A	231	ASP
1	A	235	VAL
1	A	245	LYS
1	A	250	GLU
1	A	251	THR
1	A	253	ARG
1	A	254	LEU
1	A	256	GLN
1	A	257	GLU
1	A	262	ILE
1	A	263	ILE
1	A	265	VAL
1	A	270	ILE
1	A	275	LYS
1	A	277	LEU
1	A	282	ARG
1	A	288	SER
1	A	293	LEU
1	A	297	LYS
1	A	309	ARG
1	A	310	LEU
1	A	319	VAL
1	A	324	LEU
1	A	331	CYS
1	A	337	ARG
1	A	340	LEU
1	A	354	LYS
1	A	358	LEU
1	A	369	TYR
1	A	375	LYS

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Mol	Chain	Res	Type
1	A	379	ILE
1	A	384	ARG
1	A	386	SER
1	A	394	GLU
1	A	400	ASN
1	A	409	LEU
1	A	412	ASP
1	A	417	LEU
1	A	420	ARG
1	A	434	ARG
1	A	435	LEU
1	A	443	SER
1	A	447	LEU
1	A	451	LEU
1	A	454	LEU
1	A	457	TYR
1	A	460	ILE
1	A	461	SER
1	A	466	ARG
1	A	467	SER
1	A	468	HIS
1	A	470	THR
1	A	476	LEU
1	A	497	TYR
1	A	503	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	30	GLN
1	A	159	HIS
1	A	175	GLN
1	A	336	GLN
1	A	400	ASN
1	A	418	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	PLP	A	962	1	15,15,16	1.96	6 (40%)	21,22,23	2.98	11 (52%)

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	PLP	A	962	1	-	5/6/6/8	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	962	PLP	C2-N1	3.58	1.40	1.33
2	A	962	PLP	P-O4P	-3.11	1.50	1.60
2	A	962	PLP	C6-N1	2.66	1.39	1.34
2	A	962	PLP	C5A-C5	2.55	1.57	1.50
2	A	962	PLP	C5-C4	2.49	1.43	1.40
2	A	962	PLP	P-O2P	-2.08	1.47	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	962	PLP	C5A-C5-C6	-6.30	109.08	119.36
2	A	962	PLP	C4A-C4-C5	5.71	126.83	120.94
2	A	962	PLP	C5A-C5-C4	5.63	133.75	122.64
2	A	962	PLP	O4P-C5A-C5	3.63	116.16	109.36
2	A	962	PLP	C4A-C4-C3	-3.55	114.60	120.52
2	A	962	PLP	O3-C3-C2	3.46	124.76	117.58
2	A	962	PLP	C2A-C2-N1	3.16	123.60	117.64
2	A	962	PLP	C2A-C2-C3	-3.15	117.11	120.80
2	A	962	PLP	O3-C3-C4	-2.47	111.67	118.10
2	A	962	PLP	C4-C3-C2	2.35	123.39	119.89
2	A	962	PLP	O3P-P-O2P	2.10	115.69	107.80

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	962	PLP	C4-C5-C5A-O4P
2	A	962	PLP	C6-C5-C5A-O4P
2	A	962	PLP	C5A-O4P-P-O1P
2	A	962	PLP	C5A-O4P-P-O2P
2	A	962	PLP	C5A-O4P-P-O3P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	962	PLP	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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, 95th 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^2)	Q<0.9
1	A	494/514 (96%)	-0.81	10 (2%) 65 56	8, 32, 80, 99	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	GLU	4.2
1	A	80	HIS	3.8
1	A	348	GLU	3.0
1	A	5	GLN	2.8
1	A	250	GLU	2.7
1	A	469	GLY	2.5
1	A	404	TYR	2.3
1	A	76	GLU	2.3
1	A	470	THR	2.2
1	A	457	TYR	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 monosaccharides 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, 95th 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^2)	Q<0.9
2	PLP	A	962	15/16	0.98	0.09	5,11,40,83	0

6.5 Other polymers (i)

There are no such residues in this entry.