



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:45 AM UTC

PDB ID : 1A3W / pdb_00001a3w
Title : PYRUVATE KINASE FROM SACCHAROMYCES CEREVISIAE COM-
PLEXED WITH FBP, PG, MN2+ AND K+
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Deposited on : 1998-01-26
Resolution : 3.00 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

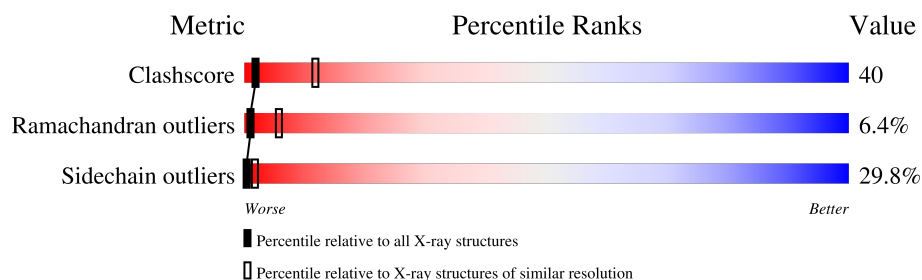
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

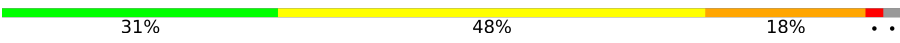
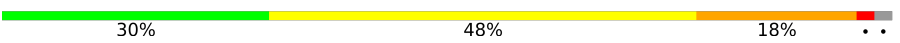
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	500	
1	B	500	

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	PGA	A	1005	-	X	-	-
2	PGA	B	1006	-	X	-	-

2 Entry composition [i](#)

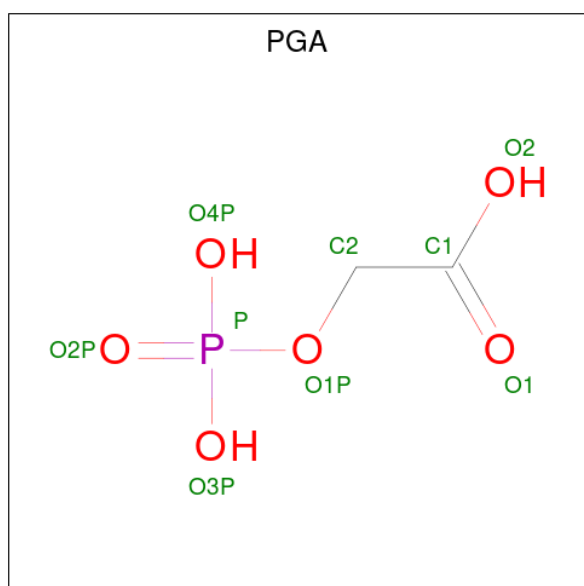
There are 5 unique types of molecules in this entry. The entry contains 7581 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PYRUVATE KINASE.

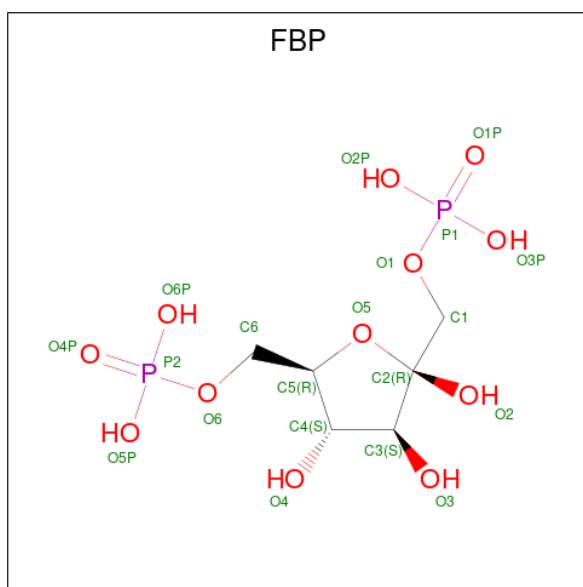
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	0	0
			3773	2380	650	726	17			
1	B	489	Total	C	N	O	S	0	0	0
			3746	2362	647	720	17			

- Molecule 2 is 2-PHOSPHOGLYCOLIC ACID (CCD ID: PGA) (formula: $C_2H_5O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			9	2	6	1		
2	B	1	Total	C	O	P	0	0
			9	2	6	1		

- Molecule 3 is 1,6-di-O-phosphono-beta-D-fructofuranose (CCD ID: FBP) (formula: $C_6H_{14}O_{12}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			20	6	12	2		
3	B	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		

- Molecule 5 is POTASSIUM ION (CCD ID: K) (formula: K).

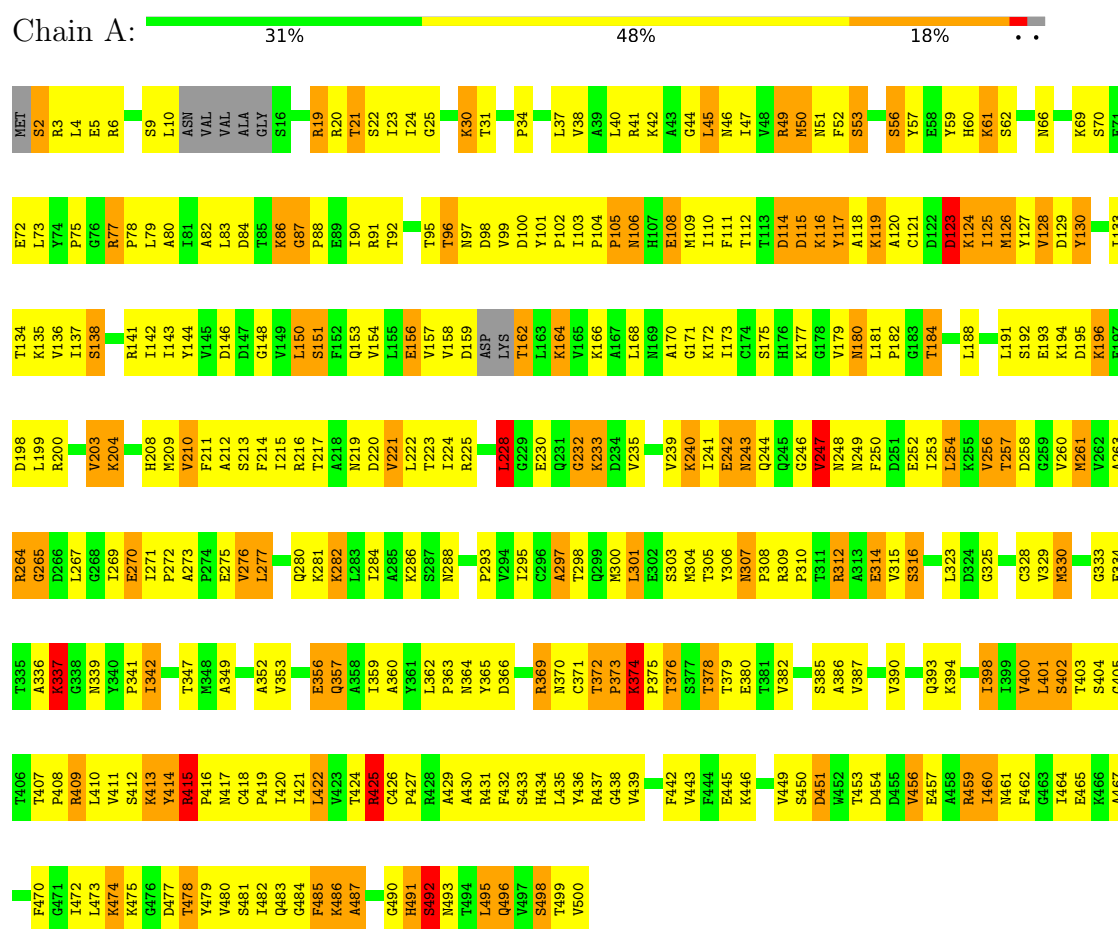
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	K	0	0
			1	1		
5	B	1	Total	K	0	0
			1	1		

3 Residue-property plots

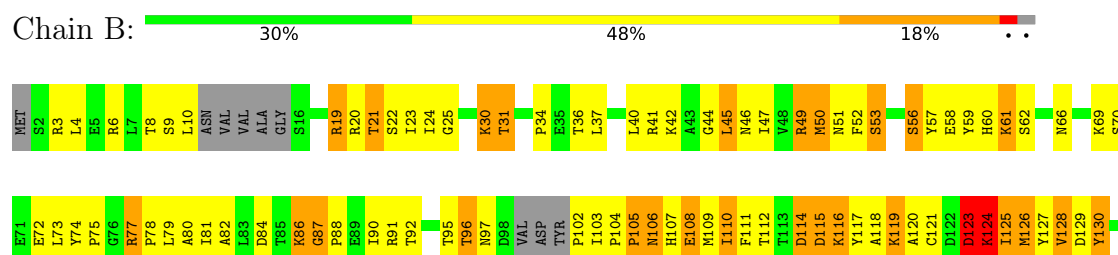
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

• Molecule 1: PYRUVATE KINASE



• Molecule 1: PYRUVATE KINASE





4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.40Å 102.70Å 110.90Å 90.00° 112.30° 90.00°	Depositor
Resolution (Å)	100.00 – 3.00	Depositor
% Data completeness (in resolution range)	75.9 (100.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.218 , 0.323	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7581	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, K, PGA, FBP

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	0.36	0/3834	0.87	7/5196 (0.1%)
1	B	0.36	0/3805	0.88	10/5153 (0.2%)
All	All	0.36	0/7639	0.87	17/10349 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	371	CYS	N-CA-C	-6.82	105.30	112.93
1	B	371	CYS	N-CA-C	-6.77	105.35	112.93
1	A	297	ALA	N-CA-C	6.38	118.47	108.96
1	B	297	ALA	N-CA-C	6.29	118.33	108.96
1	B	256	VAL	N-CA-C	6.17	116.80	110.82
1	A	372	THR	N-CA-C	-5.97	101.27	108.14
1	A	228	LEU	N-CA-C	-5.80	105.04	111.71
1	B	372	THR	N-CA-C	-5.74	101.54	108.14
1	A	247	VAL	N-CA-C	-5.72	107.27	111.90
1	A	256	VAL	N-CA-C	5.70	116.35	110.82
1	B	228	LEU	N-CA-C	-5.68	105.18	111.71
1	B	369	ARG	NE-CZ-NH2	-5.58	114.17	119.20
1	B	247	VAL	N-CA-C	-5.42	107.51	111.90
1	B	58	GLU	N-CA-C	-5.33	105.91	112.90
1	A	425	ARG	N-CA-C	-5.25	106.56	113.12
1	B	369	ARG	CD-NE-CZ	5.18	131.65	124.40
1	B	425	ARG	N-CA-C	-5.03	106.84	113.12

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3773	0	3844	307	0
1	B	3746	0	3822	313	0
2	A	9	0	2	1	0
2	B	9	0	2	0	0
3	A	20	0	10	3	0
3	B	20	0	10	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	7581	0	7690	605	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 40.

All (605) 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:374:LYS:H	1:B:375:PRO:HA	1.16	1.11
1:A:374:LYS:H	1:A:375:PRO:HA	1.19	1.05
1:B:398:ILE:HD11	1:B:482:ILE:HD11	1.43	1.01
1:A:242:GLU:HG3	1:A:263:ALA:CB	1.95	0.96
1:B:242:GLU:HG3	1:B:263:ALA:CB	1.97	0.95
1:B:242:GLU:HG3	1:B:263:ALA:HB1	1.49	0.95
1:B:301:LEU:HA	1:B:314:GLU:HG2	1.47	0.95
1:A:242:GLU:HG3	1:A:263:ALA:HB1	1.49	0.94
1:A:301:LEU:HA	1:A:314:GLU:HG2	1.48	0.94
1:A:398:ILE:HD11	1:A:482:ILE:HD11	1.48	0.92
1:B:24:ILE:HB	1:B:330:MET:HB2	1.53	0.91
1:A:10:LEU:H	1:B:282:LYS:HE2	1.37	0.90
1:A:24:ILE:HB	1:A:330:MET:HB2	1.52	0.89
1:A:119:LYS:HE2	1:A:119:LYS:H	1.40	0.86
1:A:121:CYS:HB2	1:A:126:MET:HA	1.58	0.86
1:A:282:LYS:HE2	1:B:10:LEU:H	1.40	0.86
1:B:119:LYS:HE2	1:B:119:LYS:H	1.40	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:LEU:HD22	1:A:500:VAL:HG21	1.57	0.86
1:A:310:PRO:HG3	1:A:347:THR:HG21	1.57	0.86
1:A:240:LYS:HD3	1:A:261:MET:SD	2.16	0.85
1:B:121:CYS:HB2	1:B:126:MET:HA	1.56	0.85
1:B:310:PRO:HG3	1:B:347:THR:HG21	1.58	0.84
1:A:373:PRO:O	1:A:374:LYS:HB2	1.78	0.84
1:B:473:LEU:HD22	1:B:500:VAL:HG21	1.58	0.84
1:A:373:PRO:HB2	1:A:375:PRO:HA	1.59	0.82
1:B:373:PRO:HB2	1:B:375:PRO:HA	1.61	0.82
1:B:110:ILE:HD11	1:B:162:THR:HG22	1.62	0.81
1:B:373:PRO:O	1:B:374:LYS:HB2	1.78	0.81
1:B:23:ILE:H	1:B:46:ASN:HD22	1.28	0.81
1:A:136:VAL:HG22	1:A:184:THR:HG22	1.61	0.81
1:B:136:VAL:HG22	1:B:184:THR:HG22	1.63	0.80
1:A:126:MET:HG3	1:A:173:ILE:HD11	1.63	0.79
1:A:456:VAL:HG12	1:A:459:ARG:HH21	1.46	0.79
1:B:144:TYR:HB3	1:B:148:GLY:HA2	1.65	0.79
1:A:23:ILE:H	1:A:46:ASN:HD22	1.28	0.78
1:A:151:SER:HB2	1:A:168:LEU:HD12	1.65	0.78
1:A:144:TYR:HB3	1:A:148:GLY:HA2	1.66	0.78
1:A:30:LYS:HE3	1:A:30:LYS:H	1.48	0.78
1:A:24:ILE:HB	1:A:330:MET:CB	2.13	0.78
1:B:456:VAL:HG12	1:B:459:ARG:HH21	1.48	0.78
1:B:24:ILE:HB	1:B:330:MET:CB	2.12	0.77
1:B:240:LYS:HD3	1:B:261:MET:SD	2.25	0.77
1:A:429:ALA:HA	1:A:432:PHE:CZ	2.19	0.77
1:B:30:LYS:HE3	1:B:30:LYS:H	1.48	0.77
1:B:374:LYS:H	1:B:375:PRO:CA	1.96	0.77
1:B:95:THR:HA	1:B:121:CYS:O	1.86	0.76
1:A:95:THR:HA	1:A:121:CYS:O	1.85	0.75
1:B:34:PRO:HD3	1:B:66:ASN:HB3	1.68	0.75
1:B:126:MET:HG3	1:B:173:ILE:HD11	1.67	0.75
1:A:110:ILE:HD11	1:A:162:THR:HG22	1.68	0.74
1:A:91:ARG:H	1:A:129:ASP:HB2	1.52	0.74
1:B:373:PRO:HB2	1:B:376:THR:N	2.03	0.74
1:B:429:ALA:HA	1:B:432:PHE:CZ	2.23	0.74
1:A:104:PRO:HD2	1:A:171:GLY:O	1.88	0.72
1:A:456:VAL:HG12	1:A:459:ARG:NH2	2.04	0.72
1:A:374:LYS:H	1:A:375:PRO:CA	1.99	0.72
1:A:244:GLN:O	1:A:247:VAL:HG23	1.90	0.72
1:B:91:ARG:H	1:B:129:ASP:HB2	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:PRO:HD3	1:A:66:ASN:HB3	1.71	0.72
1:A:90:ILE:HG23	1:A:130:TYR:H	1.55	0.72
1:B:301:LEU:HD23	1:B:314:GLU:HG2	1.72	0.71
1:A:301:LEU:HD23	1:A:314:GLU:HG2	1.71	0.71
1:B:241:ILE:O	1:B:263:ALA:HB3	1.91	0.71
1:A:246:GLY:O	1:A:250:PHE:HB2	1.90	0.71
1:B:386:ALA:HA	1:B:495:LEU:HD11	1.72	0.71
1:A:373:PRO:HB2	1:A:376:THR:N	2.06	0.71
1:A:413:LYS:C	1:A:415:ARG:H	1.99	0.70
1:B:115:ASP:H	1:B:127:TYR:HE2	1.39	0.70
1:A:386:ALA:HA	1:A:495:LEU:HD11	1.72	0.70
1:A:115:ASP:H	1:A:127:TYR:HE2	1.40	0.70
1:B:151:SER:HB2	1:B:168:LEU:HD12	1.73	0.70
1:B:41:ARG:HE	1:B:79:LEU:HB2	1.57	0.70
1:B:104:PRO:HD2	1:B:171:GLY:O	1.91	0.70
1:B:56:SER:N	1:B:59:TYR:HB2	2.07	0.69
1:B:456:VAL:HG12	1:B:459:ARG:NH2	2.06	0.69
1:B:374:LYS:N	1:B:375:PRO:HA	1.98	0.69
1:A:102:PRO:O	1:A:173:ILE:HG22	1.93	0.69
1:A:295:ILE:HG12	1:A:328:CYS:HB2	1.73	0.69
1:B:413:LYS:C	1:B:415:ARG:H	2.00	0.69
1:A:41:ARG:HE	1:A:79:LEU:HB2	1.58	0.68
1:A:56:SER:N	1:A:59:TYR:HB2	2.09	0.68
1:A:239:VAL:CG2	1:A:260:VAL:HG12	2.23	0.68
1:B:103:ILE:HA	1:B:172:LYS:HA	1.76	0.68
1:A:378:THR:HG21	1:A:486:LYS:HE2	1.75	0.68
1:B:373:PRO:HB2	1:B:376:THR:H	1.59	0.68
1:A:95:THR:HG23	1:A:175:SER:HB3	1.75	0.68
1:B:159:ASP:HA	1:B:162:THR:HB	1.75	0.68
1:B:246:GLY:O	1:B:250:PHE:HB2	1.94	0.68
1:A:374:LYS:N	1:A:375:PRO:HA	2.00	0.68
1:B:451:ASP:C	1:B:453:THR:H	2.02	0.68
1:A:19:ARG:HH12	1:A:46:ASN:HD21	1.42	0.67
1:B:239:VAL:CG2	1:B:260:VAL:HG12	2.23	0.67
1:B:90:ILE:HG23	1:B:130:TYR:H	1.58	0.67
1:B:432:PHE:O	1:B:435:LEU:HB2	1.95	0.67
1:B:295:ILE:HG12	1:B:328:CYS:HB2	1.76	0.67
1:A:130:TYR:HB3	1:A:133:ILE:HB	1.77	0.67
1:A:451:ASP:C	1:A:453:THR:H	2.02	0.67
1:B:19:ARG:HH12	1:B:46:ASN:HD21	1.42	0.67
1:B:378:THR:HG21	1:B:486:LYS:HE2	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ALA:O	1:B:267:LEU:HB2	1.95	0.66
1:B:398:ILE:HD11	1:B:482:ILE:CD1	2.23	0.66
1:A:199:LEU:O	1:A:203:VAL:HG13	1.95	0.66
1:A:103:ILE:HA	1:A:172:LYS:HA	1.77	0.66
1:A:360:ALA:O	1:A:364:ASN:HB2	1.96	0.66
1:B:130:TYR:HB3	1:B:133:ILE:HB	1.78	0.66
1:B:460:ILE:HD11	1:B:483:GLN:HE21	1.61	0.66
1:A:90:ILE:HD12	1:A:130:TYR:HB2	1.77	0.66
1:B:95:THR:HG23	1:B:175:SER:HB3	1.77	0.66
1:B:307:ASN:ND2	1:B:308:PRO:HD2	2.10	0.66
1:A:50:MET:HB2	1:A:82:ALA:O	1.96	0.65
1:A:96:THR:HA	1:A:120:ALA:HA	1.76	0.65
1:B:264:ARG:HH11	1:B:280:GLN:HE22	1.43	0.65
1:A:432:PHE:O	1:A:435:LEU:HB2	1.96	0.65
1:A:307:ASN:ND2	1:A:308:PRO:HD2	2.10	0.65
1:B:96:THR:HA	1:B:120:ALA:HA	1.77	0.65
1:A:110:ILE:HG13	1:A:111:PHE:N	2.12	0.65
1:B:360:ALA:O	1:B:364:ASN:HB2	1.96	0.65
1:A:209:MET:HE1	1:A:432:PHE:HA	1.78	0.65
1:B:24:ILE:HD12	1:B:295:ILE:HG21	1.77	0.65
1:B:117:TYR:HB2	1:B:127:TYR:CE1	2.32	0.65
1:B:50:MET:HB2	1:B:82:ALA:O	1.95	0.65
1:A:241:ILE:O	1:A:263:ALA:HB3	1.96	0.65
1:A:30:LYS:HE3	1:A:30:LYS:N	2.12	0.65
1:A:460:ILE:HD11	1:A:483:GLN:HE21	1.62	0.65
1:A:264:ARG:HH11	1:A:280:GLN:HE22	1.43	0.64
1:B:243:ASN:HA	1:B:270:GLU:HG2	1.78	0.64
1:A:258:ASP:O	1:A:293:PRO:HD2	1.97	0.64
1:B:30:LYS:HE3	1:B:30:LYS:N	2.11	0.64
1:A:24:ILE:HD12	1:A:295:ILE:HG21	1.79	0.64
1:A:373:PRO:HB2	1:A:376:THR:H	1.63	0.64
1:B:258:ASP:O	1:B:293:PRO:HD2	1.97	0.64
1:A:263:ALA:O	1:A:267:LEU:HB2	1.97	0.63
1:B:21:THR:HG23	1:B:437:ARG:HG3	1.81	0.63
1:A:51:ASN:OD1	1:A:53:SER:HB2	1.99	0.63
1:A:57:TYR:O	1:A:61:LYS:HB2	1.99	0.63
1:B:90:ILE:HD12	1:B:130:TYR:HB2	1.80	0.63
1:A:250:PHE:HE1	1:A:260:VAL:HG11	1.63	0.63
1:B:244:GLN:O	1:B:247:VAL:HG23	1.99	0.63
1:B:386:ALA:CA	1:B:495:LEU:HD11	2.30	0.62
1:A:484:GLY:O	1:A:485:PHE:HB2	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:TYR:O	1:B:61:LYS:HB2	1.98	0.62
1:B:159:ASP:HA	1:B:162:THR:CB	2.28	0.62
1:B:199:LEU:O	1:B:203:VAL:HG13	2.00	0.62
1:A:215:ILE:HG13	1:A:239:VAL:HG12	1.82	0.62
1:A:386:ALA:CA	1:A:495:LEU:HD11	2.30	0.62
1:B:110:ILE:HG13	1:B:111:PHE:N	2.13	0.62
1:B:209:MET:HE1	1:B:432:PHE:HA	1.81	0.61
1:A:491:HIS:HA	3:A:1007:FBP:O4	1.99	0.61
1:A:101:TYR:N	1:A:102:PRO:HD3	2.14	0.61
1:A:117:TYR:HB2	1:A:127:TYR:CE1	2.36	0.61
1:A:398:ILE:HD11	1:A:482:ILE:CD1	2.26	0.61
1:B:484:GLY:O	1:B:485:PHE:HB2	1.99	0.61
1:B:382:VAL:HG23	1:B:495:LEU:HD12	1.83	0.61
1:B:250:PHE:HE1	1:B:260:VAL:HG11	1.66	0.60
1:A:301:LEU:HB3	1:A:304:MET:HG3	1.84	0.60
1:B:47:ILE:HA	1:B:80:ALA:HB3	1.84	0.60
1:A:243:ASN:HA	1:A:270:GLU:HG2	1.82	0.60
1:A:21:THR:HG23	1:A:437:ARG:HG3	1.84	0.60
1:A:47:ILE:HA	1:A:80:ALA:HB3	1.83	0.59
1:A:239:VAL:HG23	1:A:260:VAL:HG12	1.82	0.59
1:A:244:GLN:HG3	1:A:248:ASN:ND2	2.17	0.59
1:B:244:GLN:HG3	1:B:248:ASN:ND2	2.18	0.59
1:B:239:VAL:HG23	1:B:260:VAL:HG12	1.83	0.59
1:B:53:SER:HA	1:B:86:LYS:HB2	1.84	0.59
1:A:2:SER:HB2	1:A:5:GLU:HB2	1.86	0.58
1:A:103:ILE:HG12	1:A:172:LYS:HB3	1.85	0.58
1:B:215:ILE:HG13	1:B:239:VAL:HG12	1.85	0.58
1:A:53:SER:HA	1:A:86:LYS:HB2	1.84	0.58
1:B:212:ALA:O	1:B:240:LYS:HB2	2.04	0.58
1:B:407:THR:O	1:B:411:VAL:HG23	2.03	0.58
1:B:250:PHE:CE1	1:B:260:VAL:HG11	2.39	0.58
1:B:301:LEU:HB3	1:B:304:MET:HG3	1.86	0.57
1:A:228:LEU:HG	1:A:235:VAL:HG11	1.86	0.57
1:B:103:ILE:HG12	1:B:172:LYS:HB3	1.86	0.57
1:A:250:PHE:CE1	1:A:260:VAL:HG11	2.39	0.57
1:A:316:SER:HA	1:B:277:LEU:HD12	1.86	0.57
1:A:323:LEU:HD23	1:A:359:ILE:HG12	1.86	0.57
1:A:407:THR:O	1:A:411:VAL:HG23	2.04	0.57
1:A:382:VAL:HG23	1:A:495:LEU:HD12	1.87	0.57
1:B:56:SER:H	1:B:59:TYR:HB2	1.68	0.57
1:B:117:TYR:HB2	1:B:127:TYR:HE1	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ILE:HD11	1:B:224:ILE:HD11	1.86	0.57
1:B:323:LEU:HD23	1:B:359:ILE:HG12	1.87	0.57
1:A:402:SER:HA	3:A:1007:FBP:H61	1.85	0.57
1:A:309:ARG:NH2	1:B:269:ILE:HD13	2.19	0.56
1:A:102:PRO:O	1:A:104:PRO:HD3	2.04	0.56
1:A:312:ARG:HG3	1:B:264:ARG:HB3	1.87	0.56
1:A:370:ASN:C	1:A:372:THR:H	2.13	0.56
1:A:230:GLU:CD	1:A:230:GLU:H	2.13	0.56
1:B:272:PRO:HD2	1:B:275:GLU:OE2	2.05	0.56
1:A:212:ALA:O	1:A:240:LYS:HB2	2.05	0.56
1:A:215:ILE:HD11	1:A:224:ILE:HD11	1.86	0.56
1:A:272:PRO:HD2	1:A:275:GLU:OE2	2.06	0.56
1:A:373:PRO:HB2	1:A:375:PRO:CA	2.33	0.55
1:A:100:ASP:C	1:A:102:PRO:HD3	2.31	0.55
1:B:51:ASN:OD1	1:B:53:SER:HB2	2.07	0.55
1:B:295:ILE:HG23	1:B:328:CYS:HB2	1.88	0.55
1:B:20:ARG:HH12	1:B:417:ASN:HA	1.70	0.55
1:B:307:ASN:HD22	1:B:308:PRO:HD2	1.70	0.55
1:B:370:ASN:C	1:B:372:THR:H	2.13	0.55
1:A:10:LEU:HB2	1:B:282:LYS:HG3	1.89	0.55
1:B:228:LEU:HG	1:B:235:VAL:HG11	1.89	0.55
1:A:88:PRO:HD2	1:A:214:PHE:HB2	1.88	0.55
1:B:366:ASP:O	1:B:369:ARG:HG3	2.07	0.55
1:A:56:SER:H	1:A:59:TYR:HB2	1.70	0.55
1:B:34:PRO:HD3	1:B:66:ASN:CB	2.37	0.55
1:B:373:PRO:HB2	1:B:375:PRO:CA	2.35	0.55
1:A:456:VAL:HA	1:A:459:ARG:HE	1.71	0.54
1:B:102:PRO:O	1:B:173:ILE:HG22	2.07	0.54
1:A:405:GLY:O	1:A:408:PRO:HG2	2.08	0.54
1:A:20:ARG:HH12	1:A:417:ASN:HA	1.71	0.54
1:A:307:ASN:HD22	1:A:308:PRO:HD2	1.72	0.54
1:B:24:ILE:HG23	1:B:47:ILE:HG22	1.90	0.54
1:B:230:GLU:CD	1:B:230:GLU:H	2.14	0.54
1:B:123:ASP:C	1:B:125:ILE:H	2.15	0.54
1:B:208:HIS:O	1:B:235:VAL:HG23	2.07	0.54
1:B:400:VAL:HG21	1:B:407:THR:HG22	1.89	0.54
1:A:429:ALA:HA	1:A:432:PHE:CE1	2.43	0.53
1:B:23:ILE:N	1:B:46:ASN:HD22	2.03	0.53
1:B:374:LYS:N	1:B:375:PRO:CA	2.65	0.53
1:B:288:ASN:HD21	1:B:325:GLY:HA3	1.73	0.53
1:B:352:ALA:O	1:B:356:GLU:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:459:ARG:O	1:B:462:PHE:HB3	2.09	0.53
1:A:117:TYR:HB2	1:A:127:TYR:HE1	1.73	0.53
1:B:241:ILE:HD12	1:B:246:GLY:O	2.08	0.53
1:B:456:VAL:HA	1:B:459:ARG:HE	1.72	0.53
1:A:23:ILE:N	1:A:46:ASN:HD22	2.03	0.53
1:B:221:VAL:HA	1:B:224:ILE:HD12	1.91	0.53
1:B:300:MET:O	1:B:301:LEU:HG	2.08	0.53
1:B:401:LEU:HD12	3:B:1008:FBP:O2	2.09	0.53
1:A:288:ASN:HD21	1:A:325:GLY:HA3	1.73	0.53
1:A:400:VAL:HG21	1:A:407:THR:HG22	1.91	0.53
1:A:419:PRO:HG2	1:A:472:ILE:HG23	1.91	0.53
1:A:269:ILE:HD13	1:B:309:ARG:NH2	2.24	0.52
1:A:98:ASP:O	1:A:100:ASP:N	2.42	0.52
1:A:221:VAL:HA	1:A:224:ILE:HD12	1.91	0.52
1:A:232:GLY:C	1:A:235:VAL:HG12	2.35	0.52
1:A:24:ILE:HG23	1:A:47:ILE:HG22	1.91	0.52
1:A:352:ALA:O	1:A:356:GLU:HG3	2.09	0.52
1:B:114:ASP:C	1:B:116:LYS:H	2.17	0.52
1:B:108:GLU:HA	1:B:166:LYS:HA	1.90	0.52
1:B:240:LYS:HG2	1:B:261:MET:HB3	1.91	0.52
1:A:295:ILE:HG23	1:A:328:CYS:HB2	1.91	0.52
1:B:373:PRO:CB	1:B:376:THR:H	2.22	0.52
1:A:336:ALA:O	1:A:337:LYS:HB3	2.09	0.52
1:B:424:THR:HG21	1:B:430:ALA:HB2	1.90	0.52
1:A:109:MET:HE2	1:A:111:PHE:HE1	1.73	0.52
1:A:459:ARG:O	1:A:462:PHE:HB3	2.09	0.52
1:A:46:ASN:O	1:A:79:LEU:HD22	2.09	0.52
1:A:123:ASP:C	1:A:125:ILE:H	2.18	0.52
1:A:244:GLN:C	1:A:247:VAL:HG23	2.35	0.52
1:A:265:GLY:HA3	1:A:298:THR:HG21	1.91	0.52
1:A:360:ALA:HB1	1:A:363:PRO:HG2	1.92	0.52
1:A:374:LYS:N	1:A:375:PRO:CA	2.66	0.52
1:B:336:ALA:O	1:B:337:LYS:HB3	2.10	0.52
1:A:108:GLU:HA	1:A:166:LYS:HA	1.91	0.51
1:A:424:THR:HG21	1:A:430:ALA:HB2	1.92	0.51
1:B:360:ALA:HB1	1:B:363:PRO:HG2	1.91	0.51
1:B:405:GLY:O	1:B:408:PRO:HG2	2.11	0.51
1:A:333:GLY:HA2	1:A:336:ALA:HB3	1.91	0.51
1:A:451:ASP:C	1:A:453:THR:N	2.69	0.51
1:B:393:GLN:HB3	1:B:478:THR:HG21	1.93	0.51
1:A:114:ASP:C	1:A:116:LYS:H	2.19	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:ALA:O	1:A:353:VAL:HG23	2.10	0.51
1:B:46:ASN:O	1:B:79:LEU:HD22	2.10	0.51
1:B:419:PRO:HG2	1:B:472:ILE:HG23	1.91	0.51
1:A:208:HIS:O	1:A:235:VAL:HG23	2.11	0.51
1:B:88:PRO:HD2	1:B:214:PHE:HB2	1.92	0.51
1:A:34:PRO:HD3	1:A:66:ASN:CB	2.40	0.51
1:A:52:PHE:HD2	1:A:198:ASP:OD2	1.94	0.51
1:A:240:LYS:HG2	1:A:261:MET:HB3	1.93	0.51
1:B:232:GLY:C	1:B:235:VAL:HG12	2.36	0.51
1:A:241:ILE:HD12	1:A:246:GLY:O	2.11	0.50
1:B:301:LEU:HD22	1:B:304:MET:HG3	1.93	0.50
1:A:192:SER:O	1:A:196:LYS:HD3	2.10	0.50
1:B:23:ILE:HG22	1:B:23:ILE:O	2.11	0.50
1:B:420:ILE:N	1:B:420:ILE:HD12	2.27	0.50
1:A:300:MET:O	1:A:301:LEU:HG	2.11	0.50
1:A:24:ILE:HG12	1:A:47:ILE:HB	1.93	0.50
1:A:56:SER:O	1:A:59:TYR:HB2	2.12	0.50
1:B:192:SER:O	1:B:196:LYS:HD3	2.11	0.50
1:B:356:GLU:HA	1:B:359:ILE:HG13	1.94	0.50
1:B:19:ARG:HD2	1:B:21:THR:O	2.11	0.50
1:B:265:GLY:CA	1:B:298:THR:HG21	2.42	0.50
1:A:130:TYR:O	1:A:133:ILE:HG22	2.12	0.50
1:A:301:LEU:HD23	1:A:314:GLU:HB3	1.94	0.50
1:B:41:ARG:HH21	1:B:79:LEU:HB2	1.77	0.50
1:A:356:GLU:HA	1:A:359:ILE:HG13	1.94	0.50
1:B:349:ALA:O	1:B:353:VAL:HG23	2.12	0.50
1:B:413:LYS:C	1:B:415:ARG:N	2.70	0.50
1:B:451:ASP:C	1:B:453:THR:N	2.68	0.50
1:A:121:CYS:HA	1:A:125:ILE:HG23	1.94	0.50
1:B:49:ARG:HH11	1:B:84:ASP:CG	2.19	0.50
1:B:105:PRO:O	1:B:106:ASN:HB2	2.12	0.50
1:B:265:GLY:HA3	1:B:298:THR:HG21	1.93	0.49
1:A:242:GLU:CG	1:A:263:ALA:HB1	2.33	0.49
1:A:272:PRO:HB2	1:A:275:GLU:HG3	1.93	0.49
1:A:301:LEU:HD22	1:A:310:PRO:HB3	1.94	0.49
1:A:400:VAL:O	1:A:422:LEU:HA	2.13	0.49
1:A:484:GLY:HA3	1:A:492:SER:H	1.77	0.49
1:B:130:TYR:O	1:B:133:ILE:HG22	2.11	0.49
1:B:429:ALA:HA	1:B:432:PHE:CE1	2.47	0.49
1:A:282:LYS:HG3	1:B:10:LEU:HB2	1.94	0.49
1:A:393:GLN:HB3	1:A:478:THR:HG21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ARG:HA	1:B:45:LEU:HB3	1.94	0.49
1:B:193:GLU:CD	1:B:193:GLU:H	2.21	0.49
1:B:400:VAL:O	1:B:422:LEU:HA	2.11	0.49
1:A:115:ASP:N	1:A:127:TYR:CE2	2.81	0.49
1:B:329:VAL:HG23	1:B:329:VAL:O	2.12	0.49
1:B:109:MET:HE2	1:B:111:PHE:HE1	1.78	0.49
1:B:191:LEU:HD12	1:B:223:THR:HG22	1.95	0.49
1:B:301:LEU:HD22	1:B:310:PRO:HB3	1.94	0.49
1:A:265:GLY:CA	1:A:298:THR:HG21	2.42	0.49
1:B:56:SER:O	1:B:59:TYR:HB2	2.12	0.49
1:A:329:VAL:HG23	1:A:329:VAL:O	2.13	0.49
1:B:102:PRO:HB2	1:B:173:ILE:HG23	1.94	0.49
1:A:451:ASP:OD2	1:A:454:ASP:HB2	2.12	0.49
1:B:481:SER:HB2	1:B:496:GLN:HB3	1.94	0.49
1:A:86:LYS:O	1:A:87:GLY:O	2.31	0.49
1:B:47:ILE:HG13	1:B:80:ALA:HB3	1.95	0.49
1:B:301:LEU:HD23	1:B:314:GLU:CG	2.41	0.49
1:B:25:GLY:HA3	1:B:45:LEU:HD11	1.94	0.49
1:B:416:PRO:C	1:B:418:CYS:H	2.20	0.49
1:A:115:ASP:HA	1:A:118:ALA:HB2	1.95	0.48
1:A:156:GLU:HG2	1:A:164:LYS:HE2	1.95	0.48
1:A:192:SER:H	1:A:195:ASP:HB2	1.77	0.48
1:A:263:ALA:HB1	2:A:1005:PGA:O1	2.14	0.48
1:A:413:LYS:O	1:A:415:ARG:N	2.46	0.48
1:B:30:LYS:HD2	1:B:337:LYS:HA	1.95	0.48
1:B:378:THR:O	1:B:382:VAL:HG12	2.13	0.48
1:A:215:ILE:HD12	1:A:220:ASP:HB3	1.96	0.48
1:A:264:ARG:HB3	1:B:312:ARG:HG3	1.95	0.48
1:A:421:ILE:HD11	1:A:472:ILE:HG21	1.95	0.48
1:B:333:GLY:HA2	1:B:336:ALA:HB3	1.95	0.48
1:A:41:ARG:HA	1:A:45:LEU:HB3	1.96	0.48
1:A:49:ARG:HH11	1:A:84:ASP:CG	2.21	0.48
1:A:256:VAL:HG23	1:A:257:THR:N	2.28	0.48
1:A:401:LEU:HD21	1:A:460:ILE:HG13	1.95	0.48
1:B:20:ARG:HB3	1:B:438:GLY:HA2	1.94	0.48
1:B:424:THR:CG2	1:B:430:ALA:HB2	2.43	0.48
1:B:451:ASP:OD2	1:B:454:ASP:HB2	2.12	0.48
1:B:484:GLY:HA3	1:B:492:SER:H	1.77	0.48
1:A:235:VAL:HG13	1:A:235:VAL:O	2.13	0.48
1:A:373:PRO:CB	1:A:376:THR:H	2.27	0.48
1:A:416:PRO:C	1:A:418:CYS:H	2.21	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:LEU:HD23	1:B:314:GLU:HB3	1.94	0.48
1:B:421:ILE:HD11	1:B:472:ILE:HG21	1.95	0.48
1:A:393:GLN:O	1:A:394:LYS:HB2	2.14	0.48
1:A:409:ARG:HG3	1:A:436:TYR:CZ	2.48	0.48
1:A:92:THR:HA	1:A:127:TYR:O	2.13	0.48
1:A:117:TYR:N	1:A:117:TYR:CD1	2.81	0.48
1:A:193:GLU:H	1:A:193:GLU:CD	2.22	0.48
1:A:239:VAL:O	1:A:261:MET:N	2.47	0.48
1:A:420:ILE:HD12	1:A:420:ILE:N	2.29	0.48
1:A:175:SER:O	1:A:177:LYS:HE2	2.14	0.48
1:B:52:PHE:HD2	1:B:198:ASP:OD2	1.95	0.48
1:B:425:ARG:NH2	1:B:446:LYS:O	2.47	0.48
1:B:102:PRO:HB2	1:B:173:ILE:CG2	2.43	0.48
1:B:115:ASP:HA	1:B:118:ALA:HB2	1.95	0.48
1:A:108:GLU:H	1:A:108:GLU:HG2	1.52	0.48
1:B:24:ILE:HG12	1:B:47:ILE:HB	1.94	0.48
1:B:413:LYS:O	1:B:415:ARG:N	2.47	0.48
1:A:25:GLY:HA3	1:A:45:LEU:HD11	1.95	0.47
1:A:301:LEU:HD23	1:A:314:GLU:CG	2.40	0.47
1:A:424:THR:CG2	1:A:430:ALA:HB2	2.44	0.47
1:B:49:ARG:HD2	1:B:84:ASP:HB2	1.95	0.47
1:B:432:PHE:CD1	1:B:432:PHE:C	2.92	0.47
1:A:47:ILE:HG13	1:A:80:ALA:HB3	1.96	0.47
1:A:105:PRO:O	1:A:106:ASN:HB2	2.13	0.47
1:A:416:PRO:HG2	1:A:420:ILE:HD11	1.97	0.47
1:A:474:LYS:HB2	1:A:477:ASP:OD1	2.14	0.47
1:A:100:ASP:HB3	1:A:102:PRO:HD3	1.96	0.47
1:A:281:LYS:HE2	1:B:323:LEU:HD13	1.97	0.47
1:A:432:PHE:C	1:A:432:PHE:CD1	2.92	0.47
1:B:56:SER:H	1:B:59:TYR:HD2	1.59	0.47
1:B:117:TYR:N	1:B:117:TYR:CD1	2.82	0.47
1:B:121:CYS:HA	1:B:125:ILE:HG23	1.96	0.47
1:B:192:SER:H	1:B:195:ASP:HB2	1.79	0.47
1:A:20:ARG:HB3	1:A:438:GLY:HA2	1.96	0.47
1:A:41:ARG:HH21	1:A:79:LEU:HB2	1.79	0.47
1:B:86:LYS:O	1:B:87:GLY:O	2.32	0.47
1:B:235:VAL:HG13	1:B:235:VAL:O	2.14	0.47
1:A:301:LEU:HD13	1:A:304:MET:SD	2.55	0.47
1:A:480:VAL:HA	1:A:496:GLN:O	2.15	0.47
1:A:216:ARG:C	1:A:253:ILE:HD11	2.40	0.47
1:B:244:GLN:C	1:B:247:VAL:HG23	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:PHE:O	1:B:250:PHE:CG	2.68	0.47
1:B:272:PRO:HB2	1:B:275:GLU:HG3	1.97	0.47
1:A:192:SER:N	1:A:195:ASP:HB2	2.30	0.47
1:B:192:SER:N	1:B:195:ASP:HB2	2.30	0.47
1:A:481:SER:HB2	1:A:496:GLN:HB3	1.97	0.46
1:B:141:ARG:HD2	1:B:184:THR:HG23	1.97	0.46
1:B:156:GLU:HG2	1:B:164:LYS:HE2	1.96	0.46
1:A:19:ARG:HD2	1:A:21:THR:O	2.14	0.46
1:B:115:ASP:N	1:B:127:TYR:CE2	2.80	0.46
1:B:175:SER:O	1:B:177:LYS:HE2	2.15	0.46
1:B:456:VAL:CG2	1:B:457:GLU:N	2.78	0.46
1:B:480:VAL:HA	1:B:496:GLN:O	2.16	0.46
1:A:250:PHE:O	1:A:254:LEU:HB2	2.15	0.46
1:B:380:GLU:CA	1:B:410:LEU:HD23	2.46	0.46
1:A:277:LEU:HD12	1:B:316:SER:HA	1.96	0.46
1:B:474:LYS:HB2	1:B:477:ASP:OD1	2.15	0.46
1:A:470:PHE:HB3	1:A:472:ILE:HG13	1.97	0.46
1:A:136:VAL:HG13	1:A:137:ILE:N	2.30	0.46
1:A:409:ARG:O	1:A:412:SER:N	2.49	0.46
1:B:41:ARG:HE	1:B:79:LEU:CB	2.27	0.46
1:B:92:THR:HA	1:B:127:TYR:O	2.14	0.46
1:B:200:ARG:O	1:B:204:LYS:HB2	2.16	0.46
1:B:359:ILE:HG22	1:B:360:ALA:N	2.30	0.46
1:B:401:LEU:HD21	1:B:460:ILE:HG13	1.98	0.46
1:A:425:ARG:NH2	1:A:446:LYS:O	2.46	0.46
1:B:470:PHE:HB3	1:B:472:ILE:HG13	1.98	0.46
1:A:30:LYS:HD2	1:A:337:LYS:HA	1.96	0.46
1:A:49:ARG:HD2	1:A:84:ASP:HB2	1.97	0.46
1:A:359:ILE:HG22	1:A:360:ALA:N	2.30	0.46
1:B:24:ILE:CD1	1:B:295:ILE:HD13	2.45	0.46
1:B:491:HIS:HA	3:B:1008:FBP:O4	2.15	0.46
1:B:136:VAL:HG13	1:B:137:ILE:N	2.31	0.46
1:A:264:ARG:HD2	1:A:280:GLN:NE2	2.31	0.46
1:B:401:LEU:HD22	1:B:459:ARG:HB3	1.98	0.46
1:B:216:ARG:C	1:B:253:ILE:HD11	2.41	0.45
1:A:297:ALA:CB	1:A:330:MET:HG2	2.47	0.45
1:A:159:ASP:HA	1:A:162:THR:HB	1.98	0.45
1:A:353:VAL:HG12	1:A:357:GLN:OE1	2.16	0.45
1:B:244:GLN:HG3	1:B:248:ASN:HD21	1.79	0.45
1:B:393:GLN:O	1:B:394:LYS:HB2	2.16	0.45
1:A:362:LEU:HB3	1:A:363:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:PRO:HG2	1:B:420:ILE:HD11	1.97	0.45
1:A:244:GLN:HG3	1:A:248:ASN:HD21	1.78	0.45
1:A:380:GLU:CA	1:A:410:LEU:HD23	2.47	0.45
1:B:191:LEU:HD11	1:B:224:ILE:HA	1.98	0.45
1:A:141:ARG:HD2	1:A:184:THR:HG23	1.98	0.45
1:B:102:PRO:O	1:B:104:PRO:HD3	2.17	0.45
1:B:264:ARG:HD2	1:B:280:GLN:NE2	2.31	0.45
1:B:49:ARG:HG2	1:B:211:PHE:CE2	2.52	0.45
1:B:241:ILE:HG23	1:B:246:GLY:C	2.42	0.45
1:B:256:VAL:HG23	1:B:257:THR:N	2.32	0.45
1:B:293:PRO:HG3	1:B:436:TYR:HE1	1.81	0.45
1:A:200:ARG:O	1:A:204:LYS:HB2	2.16	0.45
1:B:239:VAL:O	1:B:261:MET:N	2.46	0.45
1:B:74:TYR:HA	1:B:75:PRO:HD2	1.71	0.45
1:B:114:ASP:CG	1:B:116:LYS:HB2	2.41	0.45
1:B:250:PHE:O	1:B:254:LEU:HB2	2.17	0.45
1:B:362:LEU:HB3	1:B:363:PRO:HD3	1.99	0.45
1:A:56:SER:H	1:A:59:TYR:HD2	1.60	0.44
1:A:242:GLU:HG3	1:A:263:ALA:HB3	1.91	0.44
1:B:409:ARG:O	1:B:412:SER:N	2.49	0.44
1:A:24:ILE:CD1	1:A:295:ILE:HD13	2.47	0.44
1:A:79:LEU:HD13	1:A:79:LEU:C	2.43	0.44
1:A:138:SER:H	1:A:141:ARG:HG3	1.82	0.44
1:A:456:VAL:CG2	1:A:457:GLU:N	2.80	0.44
1:B:24:ILE:HD11	1:B:295:ILE:HD13	1.99	0.44
1:A:191:LEU:HD11	1:A:224:ILE:HA	1.99	0.44
1:A:312:ARG:NH1	1:B:268:GLY:HA3	2.32	0.44
1:B:40:LEU:HD12	1:B:342:ILE:HG22	1.99	0.44
1:B:77:ARG:NH2	1:B:442:PHE:CE1	2.85	0.44
1:B:460:ILE:O	1:B:464:ILE:HG13	2.18	0.44
1:B:242:GLU:CG	1:B:263:ALA:HB1	2.33	0.44
1:B:356:GLU:HG3	1:B:356:GLU:H	1.63	0.44
1:A:77:ARG:NH2	1:A:442:PHE:CE1	2.86	0.44
1:A:128:VAL:HG12	1:A:179:VAL:HG21	2.00	0.44
1:B:114:ASP:OD1	1:B:116:LYS:HB2	2.18	0.44
1:B:239:VAL:HG23	1:B:260:VAL:HA	2.00	0.44
1:A:362:LEU:HB3	1:A:363:PRO:CD	2.48	0.44
1:A:412:SER:O	1:A:415:ARG:HG3	2.18	0.44
1:B:49:ARG:HG2	1:B:211:PHE:HE2	1.83	0.44
1:B:138:SER:H	1:B:141:ARG:HG3	1.82	0.44
1:A:133:ILE:HA	1:A:136:VAL:HG12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:TYR:HA	1:A:150:LEU:O	2.18	0.43
1:A:286:LYS:HG2	1:B:8:THR:HG23	2.00	0.43
1:B:133:ILE:O	1:B:137:ILE:HG22	2.18	0.43
1:B:142:ILE:C	1:B:143:ILE:HG13	2.43	0.43
1:B:234:ASP:OD1	1:B:428:ARG:NH1	2.51	0.43
1:B:353:VAL:HG12	1:B:357:GLN:OE1	2.18	0.43
1:B:115:ASP:N	1:B:127:TYR:OH	2.51	0.43
1:A:456:VAL:CG1	1:A:459:ARG:HH21	2.24	0.43
1:A:40:LEU:HD12	1:A:342:ILE:HG22	1.99	0.43
1:A:49:ARG:HG2	1:A:211:PHE:CE2	2.53	0.43
1:A:239:VAL:HG23	1:A:260:VAL:HA	2.00	0.43
1:A:142:ILE:C	1:A:143:ILE:HG13	2.44	0.43
1:A:216:ARG:O	1:A:253:ILE:HD11	2.19	0.43
1:B:242:GLU:HG2	1:B:266:ASP:HB2	2.01	0.43
1:B:273:ALA:O	1:B:276:VAL:HG12	2.19	0.43
1:B:467:ALA:CB	1:B:473:LEU:HD12	2.47	0.43
1:A:133:ILE:O	1:A:137:ILE:HG22	2.18	0.43
1:A:228:LEU:HG	1:A:235:VAL:CG1	2.48	0.43
1:B:114:ASP:OD2	1:B:116:LYS:HB2	2.19	0.43
1:B:363:PRO:O	1:B:366:ASP:HB2	2.18	0.43
1:A:217:THR:O	1:A:220:ASP:HB2	2.19	0.43
1:A:301:LEU:HD22	1:A:304:MET:HG3	2.01	0.43
1:A:34:PRO:O	1:A:38:VAL:N	2.51	0.43
1:A:90:ILE:HB	1:A:179:VAL:HB	2.01	0.43
1:A:180:ASN:C	1:A:182:PRO:HD3	2.44	0.43
1:B:79:LEU:C	1:B:79:LEU:HD13	2.43	0.43
1:B:243:ASN:HB2	1:B:244:GLN:H	1.72	0.43
1:B:301:LEU:HD13	1:B:304:MET:SD	2.59	0.43
1:A:191:LEU:HD12	1:A:223:THR:HG22	2.01	0.43
1:A:366:ASP:O	1:A:369:ARG:HG3	2.19	0.43
1:B:207:VAL:C	1:B:428:ARG:HH22	2.27	0.43
1:B:215:ILE:HD12	1:B:220:ASP:HB3	2.01	0.43
1:B:264:ARG:O	1:B:265:GLY:C	2.61	0.43
1:B:297:ALA:CB	1:B:330:MET:HG2	2.48	0.43
1:B:373:PRO:HB2	1:B:375:PRO:C	2.43	0.43
1:B:456:VAL:CG1	1:B:459:ARG:HH21	2.26	0.43
1:B:460:ILE:HD11	1:B:483:GLN:NE2	2.31	0.43
1:A:264:ARG:O	1:A:265:GLY:C	2.61	0.42
1:A:387:VAL:HG11	1:A:414:TYR:O	2.19	0.42
1:A:401:LEU:HD22	1:A:459:ARG:HB3	2.00	0.42
1:B:23:ILE:H	1:B:46:ASN:ND2	2.07	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:PHE:O	1:A:250:PHE:CG	2.71	0.42
1:A:433:SER:HB3	1:A:439:VAL:HG11	2.01	0.42
1:A:273:ALA:O	1:A:276:VAL:HG12	2.20	0.42
1:A:363:PRO:O	1:A:366:ASP:HB2	2.19	0.42
1:B:81:ILE:N	1:B:208:HIS:ND1	2.58	0.42
1:A:293:PRO:HG3	1:A:436:TYR:HE1	1.85	0.42
1:B:380:GLU:HA	1:B:410:LEU:HD23	2.01	0.42
1:B:387:VAL:HG11	1:B:414:TYR:O	2.19	0.42
1:A:373:PRO:HB2	1:A:375:PRO:C	2.43	0.42
1:B:180:ASN:C	1:B:182:PRO:HD3	2.44	0.42
1:B:303:SER:HB3	1:B:314:GLU:OE1	2.20	0.42
1:A:41:ARG:HE	1:A:79:LEU:CB	2.28	0.42
1:A:115:ASP:N	1:A:127:TYR:OH	2.52	0.42
1:A:257:THR:OG1	1:A:258:ASP:N	2.53	0.42
1:A:304:MET:HA	1:A:307:ASN:O	2.20	0.42
1:A:365:TYR:CD1	1:A:414:TYR:HB3	2.55	0.42
1:A:378:THR:O	1:A:382:VAL:HG12	2.20	0.42
1:A:460:ILE:O	1:A:464:ILE:HG13	2.19	0.42
1:B:41:ARG:HE	1:B:79:LEU:CG	2.32	0.42
1:B:114:ASP:O	1:B:116:LYS:N	2.52	0.42
1:A:23:ILE:HG22	1:A:23:ILE:O	2.19	0.42
1:B:295:ILE:HG23	1:B:328:CYS:CB	2.50	0.42
1:A:24:ILE:HD11	1:A:295:ILE:HD13	2.02	0.42
1:A:309:ARG:HH21	1:B:269:ILE:HD13	1.82	0.42
1:A:467:ALA:CB	1:A:473:LEU:HD12	2.49	0.42
1:B:107:HIS:ND1	1:B:108:GLU:O	2.47	0.42
1:B:464:ILE:O	1:B:468:LYS:HG3	2.19	0.42
1:A:117:TYR:CD2	1:A:125:ILE:HD13	2.54	0.42
1:A:380:GLU:HA	1:A:410:LEU:HD23	2.02	0.42
1:A:456:VAL:O	1:A:460:ILE:HD12	2.20	0.42
1:A:60:HIS:C	1:A:62:SER:N	2.77	0.41
1:B:194:LYS:O	1:B:198:ASP:OD1	2.38	0.41
1:A:83:LEU:HB3	1:A:210:VAL:HG22	2.02	0.41
1:A:422:LEU:O	1:A:422:LEU:HG	2.18	0.41
1:A:486:LYS:H	1:A:490:GLY:HA2	1.84	0.41
1:B:90:ILE:HB	1:B:179:VAL:HB	2.01	0.41
1:B:241:ILE:HG23	1:B:246:GLY:O	2.20	0.41
1:A:486:LYS:HB2	1:A:490:GLY:HA2	2.01	0.41
1:B:159:ASP:HA	1:B:162:THR:OG1	2.20	0.41
1:B:222:LEU:O	1:B:225:ARG:HB3	2.20	0.41
1:A:114:ASP:CG	1:A:116:LYS:HB2	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ASP:O	1:A:116:LYS:N	2.53	0.41
1:B:31:THR:HA	1:B:36:THR:HG21	2.03	0.41
1:B:110:ILE:HG23	1:B:125:ILE:HG13	2.02	0.41
1:B:144:TYR:HA	1:B:150:LEU:O	2.20	0.41
1:B:201:PHE:CD1	1:B:201:PHE:C	2.99	0.41
1:B:216:ARG:O	1:B:253:ILE:HD11	2.21	0.41
1:B:470:PHE:C	1:B:472:ILE:H	2.28	0.41
1:A:470:PHE:C	1:A:472:ILE:H	2.28	0.41
1:B:218:ALA:O	1:B:221:VAL:HG23	2.20	0.41
1:B:305:THR:HG23	1:B:306:TYR:CD2	2.55	0.41
1:B:433:SER:HB3	1:B:439:VAL:HG11	2.01	0.41
1:A:88:PRO:HB2	1:A:188:LEU:HB3	2.03	0.41
1:A:106:ASN:HD21	1:A:166:LYS:NZ	2.17	0.41
1:A:241:ILE:HG23	1:A:246:GLY:C	2.46	0.41
1:B:128:VAL:HG12	1:B:179:VAL:HG21	2.01	0.41
1:A:323:LEU:HD13	1:B:281:LYS:HE2	2.02	0.41
1:B:60:HIS:C	1:B:62:SER:N	2.78	0.41
1:B:486:LYS:HB2	1:B:490:GLY:HA2	2.02	0.41
1:A:110:ILE:HG23	1:A:125:ILE:HG13	2.01	0.41
1:A:106:ASN:ND2	1:A:106:ASN:O	2.54	0.41
1:A:225:ARG:NH2	1:A:258:ASP:OD2	2.44	0.41
1:A:264:ARG:O	1:A:267:LEU:N	2.54	0.41
1:A:486:LYS:O	1:A:487:ALA:C	2.64	0.41
1:B:56:SER:O	1:B:59:TYR:N	2.54	0.41
1:B:117:TYR:CD2	1:B:125:ILE:HD13	2.56	0.41
1:B:124:LYS:O	1:B:125:ILE:HB	2.20	0.41
1:B:124:LYS:H	1:B:124:LYS:HG2	1.70	0.41
1:B:217:THR:O	1:B:220:ASP:HB2	2.21	0.41
1:B:362:LEU:HB3	1:B:363:PRO:CD	2.51	0.41
1:B:365:TYR:CD1	1:B:414:TYR:HB3	2.56	0.41
1:B:373:PRO:CB	1:B:376:THR:N	2.79	0.41
1:B:422:LEU:O	1:B:422:LEU:HG	2.21	0.41
1:B:432:PHE:C	1:B:432:PHE:HD1	2.28	0.41
1:B:412:SER:O	1:B:415:ARG:HG3	2.21	0.41
1:A:390:VAL:O	1:A:394:LYS:N	2.54	0.40
1:A:426:CYS:HA	1:A:427:PRO:HD3	1.94	0.40
1:B:237:ILE:C	1:B:238:ILE:HD13	2.46	0.40
1:A:83:LEU:C	1:A:83:LEU:HD23	2.45	0.40
1:A:401:LEU:HD12	3:A:1007:FBP:O2	2.21	0.40
1:B:133:ILE:HA	1:B:136:VAL:HG12	2.02	0.40
1:B:466:LYS:O	1:B:469:GLU:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:LYS:O	1:B:487:ALA:C	2.63	0.40
1:A:47:ILE:HD11	1:A:434:HIS:HB2	2.03	0.40
1:B:108:GLU:H	1:B:108:GLU:HG2	1.51	0.40
1:A:432:PHE:C	1:A:432:PHE:HD1	2.29	0.40
1:B:123:ASP:C	1:B:125:ILE:N	2.80	0.40
1:A:41:ARG:HE	1:A:79:LEU:CG	2.35	0.40
1:A:49:ARG:HG2	1:A:211:PHE:HE2	1.85	0.40
1:A:114:ASP:OD1	1:A:116:LYS:HB2	2.22	0.40
1:A:116:LYS:HB3	1:A:116:LYS:NZ	2.36	0.40
1:A:305:THR:HG23	1:A:306:TYR:CD2	2.56	0.40
1:A:477:ASP:O	1:A:499:THR:HG23	2.21	0.40
1:A:479:TYR:CE1	1:A:498:SER:HB3	2.56	0.40
1:B:19:ARG:HG3	1:B:352:ALA:HB3	2.03	0.40
1:B:486:LYS:H	1:B:490:GLY:HA2	1.85	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	486/500 (97%)	398 (82%)	56 (12%)	32 (7%)	1	5
1	B	481/500 (96%)	394 (82%)	57 (12%)	30 (6%)	1	6
All	All	967/1000 (97%)	792 (82%)	113 (12%)	62 (6%)	1	6

All (62) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	78	PRO
1	A	99	VAL
1	A	106	ASN
1	A	170	ALA

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Mol	Chain	Res	Type
1	A	337	LYS
1	A	373	PRO
1	A	374	LYS
1	A	449	VAL
1	A	475	LYS
1	A	485	PHE
1	B	78	PRO
1	B	106	ASN
1	B	170	ALA
1	B	337	LYS
1	B	373	PRO
1	B	374	LYS
1	B	449	VAL
1	B	475	LYS
1	B	485	PHE
1	A	87	GLY
1	A	123	ASP
1	A	130	TYR
1	A	232	GLY
1	A	233	LYS
1	A	415	ARG
1	A	487	ALA
1	A	492	SER
1	A	493	ASN
1	B	87	GLY
1	B	123	ASP
1	B	130	TYR
1	B	154	VAL
1	B	232	GLY
1	B	233	LYS
1	B	415	ARG
1	B	487	ALA
1	B	492	SER
1	B	493	ASN
1	A	124	LYS
1	A	154	VAL
1	A	341	PRO
1	A	414	TYR
1	B	124	LYS
1	B	341	PRO
1	B	414	TYR
1	A	115	ASP

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Mol	Chain	Res	Type
1	A	146	ASP
1	B	115	ASP
1	B	146	ASP
1	B	257	THR
1	A	257	THR
1	B	301	LEU
1	A	301	LEU
1	A	105	PRO
1	B	105	PRO
1	A	44	GLY
1	A	125	ILE
1	A	265	GLY
1	B	44	GLY
1	B	125	ILE
1	B	265	GLY
1	A	75	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	417/423 (99%)	293 (70%)	124 (30%)	0	2
1	B	414/423 (98%)	290 (70%)	124 (30%)	0	1
All	All	831/846 (98%)	583 (70%)	248 (30%)	0	2

All (248) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	3	ARG
1	A	4	LEU
1	A	6	ARG
1	A	9	SER
1	A	19	ARG
1	A	21	THR

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Mol	Chain	Res	Type
1	A	22	SER
1	A	30	LYS
1	A	31	THR
1	A	37	LEU
1	A	42	LYS
1	A	45	LEU
1	A	49	ARG
1	A	50	MET
1	A	53	SER
1	A	56	SER
1	A	61	LYS
1	A	69	LYS
1	A	70	SER
1	A	72	GLU
1	A	73	LEU
1	A	77	ARG
1	A	86	LYS
1	A	96	THR
1	A	97	ASN
1	A	108	GLU
1	A	112	THR
1	A	114	ASP
1	A	116	LYS
1	A	117	TYR
1	A	119	LYS
1	A	123	ASP
1	A	124	LYS
1	A	126	MET
1	A	128	VAL
1	A	134	THR
1	A	135	LYS
1	A	138	SER
1	A	150	LEU
1	A	151	SER
1	A	153	GLN
1	A	156	GLU
1	A	157	VAL
1	A	158	VAL
1	A	162	THR
1	A	164	LYS
1	A	180	ASN
1	A	181	LEU

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Mol	Chain	Res	Type
1	A	184	THR
1	A	194	LYS
1	A	196	LYS
1	A	203	VAL
1	A	204	LYS
1	A	210	VAL
1	A	213	SER
1	A	219	ASN
1	A	221	VAL
1	A	222	LEU
1	A	228	LEU
1	A	233	LYS
1	A	240	LYS
1	A	242	GLU
1	A	243	ASN
1	A	247	VAL
1	A	249	ASN
1	A	252	GLU
1	A	254	LEU
1	A	261	MET
1	A	264	ARG
1	A	270	GLU
1	A	271	ILE
1	A	276	VAL
1	A	277	LEU
1	A	282	LYS
1	A	284	ILE
1	A	303	SER
1	A	307	ASN
1	A	312	ARG
1	A	314	GLU
1	A	315	VAL
1	A	316	SER
1	A	330	MET
1	A	334	GLU
1	A	337	LYS
1	A	339	ASN
1	A	342	ILE
1	A	356	GLU
1	A	357	GLN
1	A	369	ARG
1	A	374	LYS

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Mol	Chain	Res	Type
1	A	376	THR
1	A	378	THR
1	A	379	THR
1	A	385	SER
1	A	398	ILE
1	A	400	VAL
1	A	401	LEU
1	A	402	SER
1	A	403	THR
1	A	404	SER
1	A	409	ARG
1	A	413	LYS
1	A	415	ARG
1	A	422	LEU
1	A	425	ARG
1	A	431	ARG
1	A	443	VAL
1	A	445	GLU
1	A	450	SER
1	A	451	ASP
1	A	456	VAL
1	A	459	ARG
1	A	460	ILE
1	A	461	ASN
1	A	465	GLU
1	A	474	LYS
1	A	478	THR
1	A	486	LYS
1	A	491	HIS
1	A	492	SER
1	A	495	LEU
1	A	496	GLN
1	A	498	SER
1	B	3	ARG
1	B	4	LEU
1	B	6	ARG
1	B	9	SER
1	B	19	ARG
1	B	21	THR
1	B	22	SER
1	B	30	LYS
1	B	31	THR

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Mol	Chain	Res	Type
1	B	37	LEU
1	B	42	LYS
1	B	45	LEU
1	B	49	ARG
1	B	50	MET
1	B	53	SER
1	B	56	SER
1	B	61	LYS
1	B	69	LYS
1	B	70	SER
1	B	72	GLU
1	B	73	LEU
1	B	77	ARG
1	B	86	LYS
1	B	96	THR
1	B	97	ASN
1	B	108	GLU
1	B	110	ILE
1	B	112	THR
1	B	114	ASP
1	B	116	LYS
1	B	119	LYS
1	B	123	ASP
1	B	124	LYS
1	B	126	MET
1	B	128	VAL
1	B	134	THR
1	B	135	LYS
1	B	138	SER
1	B	150	LEU
1	B	151	SER
1	B	153	GLN
1	B	156	GLU
1	B	157	VAL
1	B	158	VAL
1	B	162	THR
1	B	164	LYS
1	B	173	ILE
1	B	180	ASN
1	B	181	LEU
1	B	184	THR
1	B	194	LYS

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Mol	Chain	Res	Type
1	B	196	LYS
1	B	203	VAL
1	B	204	LYS
1	B	210	VAL
1	B	213	SER
1	B	219	ASN
1	B	221	VAL
1	B	222	LEU
1	B	228	LEU
1	B	233	LYS
1	B	240	LYS
1	B	242	GLU
1	B	243	ASN
1	B	247	VAL
1	B	249	ASN
1	B	252	GLU
1	B	254	LEU
1	B	261	MET
1	B	264	ARG
1	B	270	GLU
1	B	271	ILE
1	B	276	VAL
1	B	277	LEU
1	B	282	LYS
1	B	283	LEU
1	B	284	ILE
1	B	303	SER
1	B	305	THR
1	B	307	ASN
1	B	312	ARG
1	B	314	GLU
1	B	315	VAL
1	B	316	SER
1	B	330	MET
1	B	337	LYS
1	B	339	ASN
1	B	342	ILE
1	B	356	GLU
1	B	357	GLN
1	B	369	ARG
1	B	374	LYS
1	B	376	THR

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Mol	Chain	Res	Type
1	B	379	THR
1	B	382	VAL
1	B	385	SER
1	B	398	ILE
1	B	400	VAL
1	B	401	LEU
1	B	402	SER
1	B	403	THR
1	B	404	SER
1	B	413	LYS
1	B	415	ARG
1	B	422	LEU
1	B	425	ARG
1	B	431	ARG
1	B	443	VAL
1	B	445	GLU
1	B	450	SER
1	B	451	ASP
1	B	456	VAL
1	B	459	ARG
1	B	460	ILE
1	B	461	ASN
1	B	465	GLU
1	B	474	LYS
1	B	478	THR
1	B	486	LYS
1	B	491	HIS
1	B	492	SER
1	B	495	LEU
1	B	496	GLN
1	B	498	SER

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

Mol	Chain	Res	Type
1	A	46	ASN
1	A	97	ASN
1	A	106	ASN
1	A	231	GLN
1	A	280	GLN
1	A	288	ASN
1	A	307	ASN

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Mol	Chain	Res	Type
1	A	320	ASN
1	A	393	GLN
1	A	434	HIS
1	A	483	GLN
1	B	46	ASN
1	B	97	ASN
1	B	106	ASN
1	B	231	GLN
1	B	280	GLN
1	B	288	ASN
1	B	307	ASN
1	B	320	ASN
1	B	364	ASN
1	B	393	GLN
1	B	434	HIS
1	B	483	GLN
1	B	491	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
3	FBP	A	1007	-	18,20,20	1.04	0	21,32,32	0.66	0
3	FBP	B	1008	-	18,20,20	1.07	1 (5%)	21,32,32	0.73	0
2	PGA	A	1005	4,5	8,8,8	2.86	4 (50%)	10,11,11	3.45	6 (60%)
2	PGA	B	1006	4,5	8,8,8	2.45	3 (37%)	10,11,11	3.45	6 (60%)

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
3	FBP	A	1007	-	-	2/13/32/32	0/1/1/1
3	FBP	B	1008	-	-	1/13/32/32	0/1/1/1
2	PGA	A	1005	4,5	-	0/6/6/6	-
2	PGA	B	1006	4,5	-	1/6/6/6	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1005	PGA	O1P-C2	-5.90	1.37	1.43
2	B	1006	PGA	P-O3P	4.33	1.70	1.54
2	A	1005	PGA	P-O3P	4.22	1.70	1.54
2	B	1006	PGA	O1P-C2	-3.91	1.39	1.43
2	B	1006	PGA	P-O2P	2.68	1.58	1.50
2	A	1005	PGA	P-O2P	2.58	1.58	1.50
3	B	1008	FBP	O5-C2	2.14	1.46	1.43
2	A	1005	PGA	P-O4P	-2.06	1.47	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1005	PGA	O1P-P-O2P	8.05	128.19	106.44
2	B	1006	PGA	O1P-P-O2P	7.59	126.95	106.44
2	B	1006	PGA	O1P-C2-C1	4.91	117.94	110.54
2	A	1005	PGA	O1P-C2-C1	4.31	117.04	110.54
2	B	1006	PGA	O2-C1-O1	-3.23	115.02	123.33
2	B	1006	PGA	O3P-P-O1P	3.17	114.94	106.67
2	A	1005	PGA	O3P-P-O1P	3.06	114.64	106.67
2	A	1005	PGA	O2-C1-O1	-2.92	115.83	123.33
2	A	1005	PGA	O4P-P-O1P	-2.43	100.32	106.67
2	B	1006	PGA	O4P-P-O1P	-2.12	101.13	106.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1006	PGA	O4P-P-O3P	-2.09	99.95	107.80
2	A	1005	PGA	O4P-P-O2P	-2.03	102.92	110.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

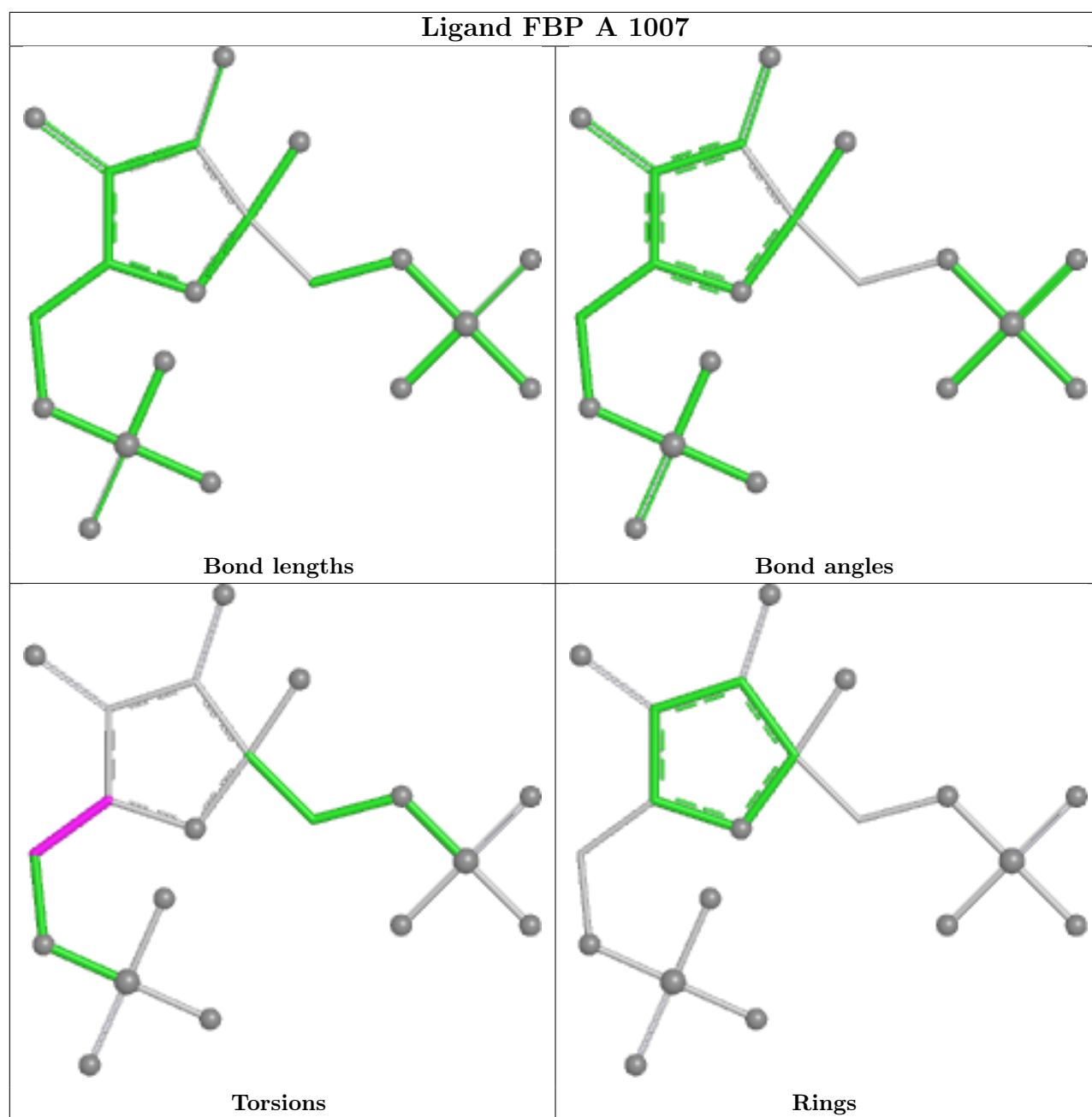
Mol	Chain	Res	Type	Atoms
3	A	1007	FBP	C4-C5-C6-O6
3	A	1007	FBP	O5-C5-C6-O6
3	B	1008	FBP	C4-C5-C6-O6
2	B	1006	PGA	C2-O1P-P-O3P

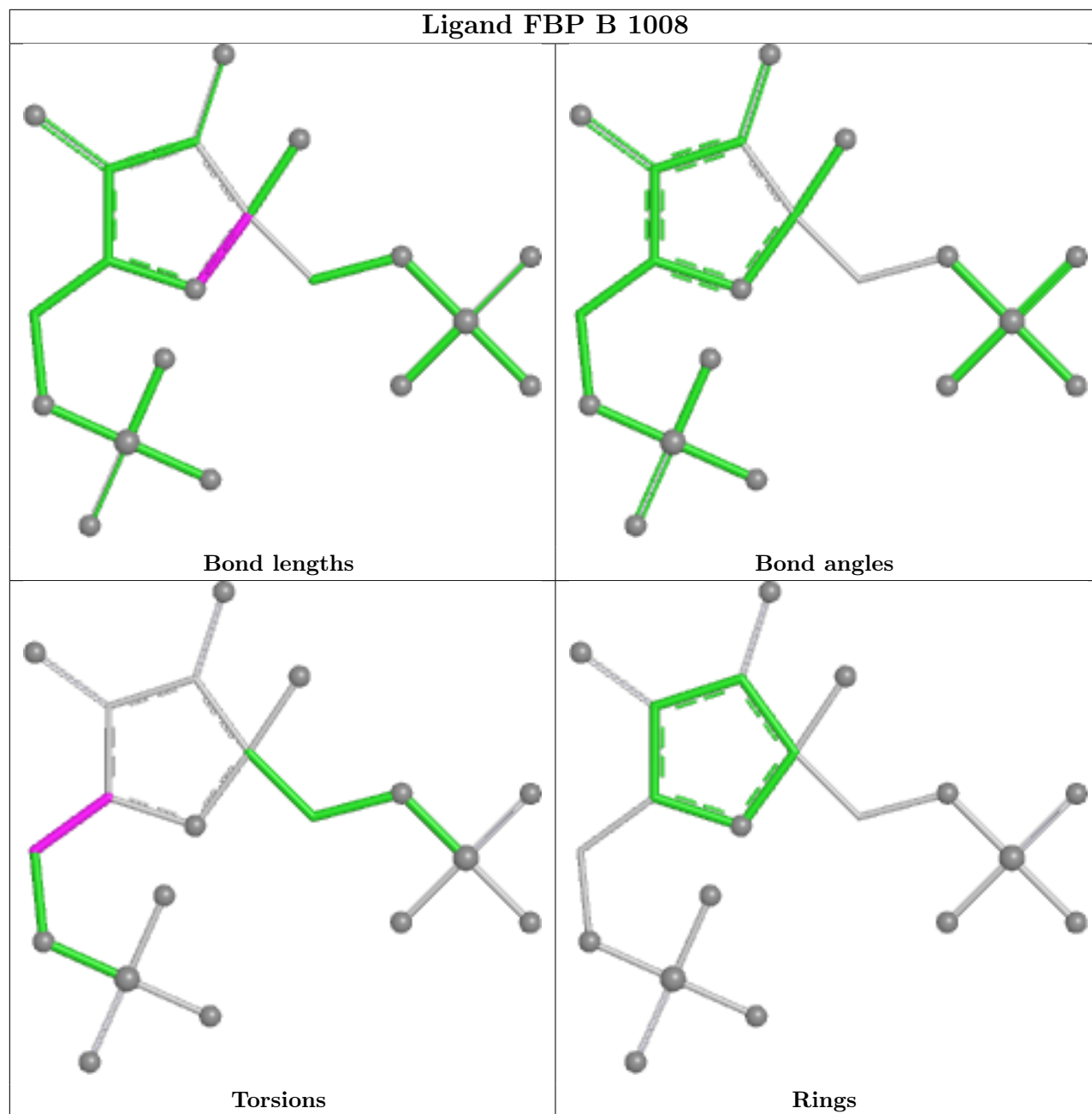
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1007	FBP	3	0
3	B	1008	FBP	2	0
2	A	1005	PGA	1	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.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.