



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:06 PM UTC

PDB ID : 2BHL / pdb_00002bhl
Title : X-RAY STRUCTURE OF HUMAN GLUCOSE-6-PHOSPHATE DEHYDROGENASE (DELETION VARIANT) COMPLEXED WITH GLUCOSE-6-PHOSPHATE
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Deposited on : 2005-01-13
Resolution : 2.90 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

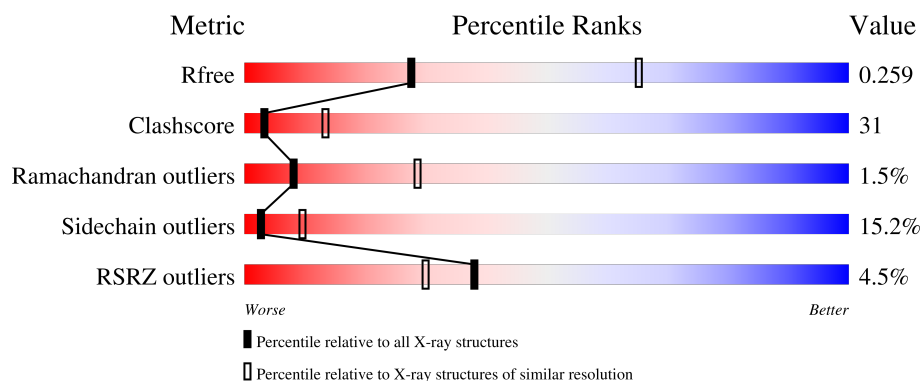
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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 $\leq 5\%$. 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	489	5% 43% 46% 8%
1	B	489	4% 46% 43% 8%

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	GOL	A	1507	-	-	X	-
2	GOL	B	1506	-	-	X	-

2 Entry composition [i](#)

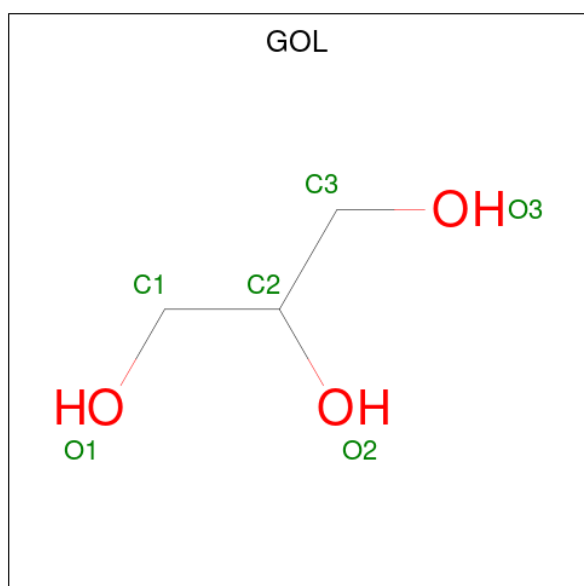
There are 4 unique types of molecules in this entry. The entry contains 7903 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLUCOSE-6-PHOSPHATE 1-DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	0	0
			3875	2471	674	710	20			
1	B	479	Total	C	N	O	S	0	0	0
			3879	2473	674	712	20			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



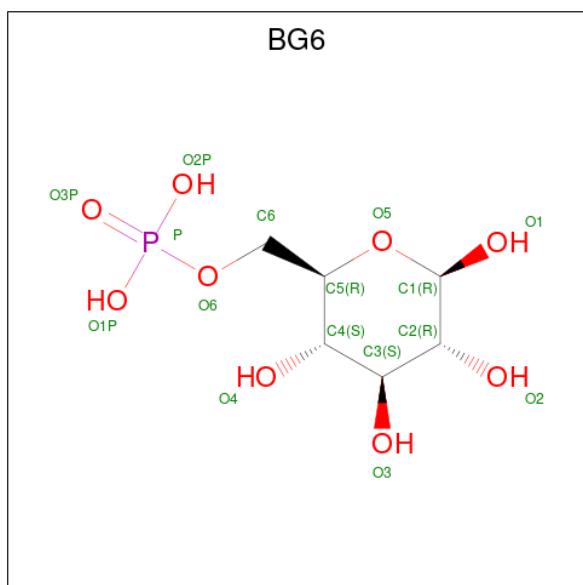
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is 6-O-phosphono-beta-D-glucopyranose (CCD ID: BG6) (formula: $C_6H_{13}O_9P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		

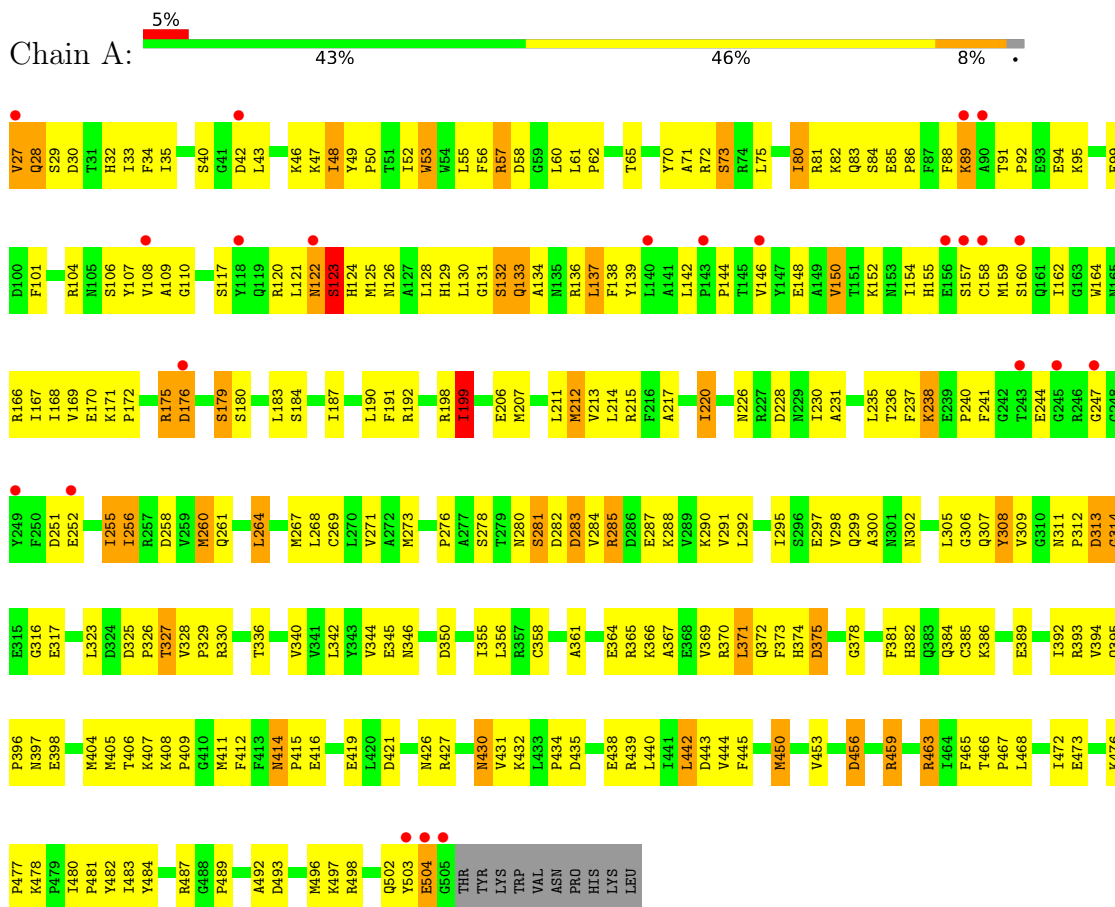
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	43	Total	O	0	0
			43	43		
4	B	38	Total	O	0	0
			38	38		

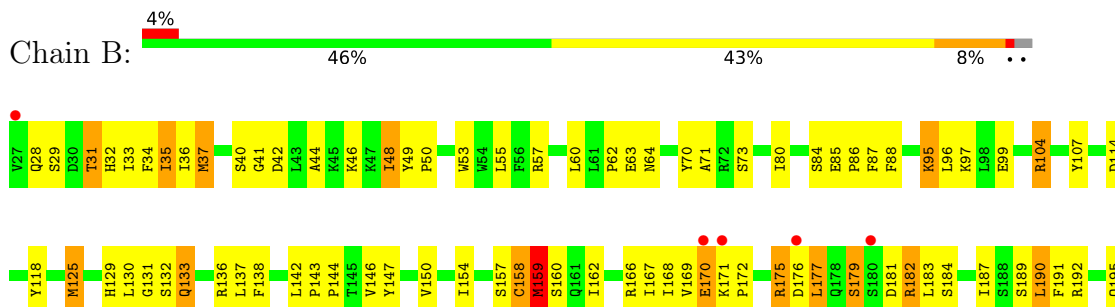
3 Residue-property plots

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 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: GLUCOSE-6-PHOSPHATE 1-DEHYDROGENASE



• Molecule 1: GLUCOSE-6-PHOSPHATE 1-DEHYDROGENASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	117.55Å 179.53Å 137.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	69.00 – 2.90 69.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (69.00-2.90) 98.7 (69.00-2.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.91Å)	Xtriage
Refinement program	TNT BUSTER/TNT	Depositor
R, R_{free}	0.212 , 0.261 0.209 , 0.259	Depositor DCC
R_{free} test set	1640 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	53.0	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 63.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7903	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, BG6

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	0.46	0/3966	0.99	15/5363 (0.3%)
1	B	0.48	0/3970	1.00	18/5368 (0.3%)
All	All	0.47	0/7936	0.99	33/10731 (0.3%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	199	ILE	N-CA-C	8.56	120.90	108.58
1	A	283	ASP	N-CA-C	-7.99	103.67	113.41
1	B	504	GLU	N-CA-C	7.56	126.89	110.80
1	A	199	ILE	N-CA-C	7.40	119.26	108.53
1	B	328	VAL	N-CA-C	6.90	113.84	107.56
1	A	327	THR	N-CA-C	-6.65	104.42	112.54
1	B	159	MET	N-CA-C	6.54	119.90	109.24
1	B	308	TYR	N-CA-C	6.47	118.82	108.41
1	A	134	ALA	N-CA-C	6.43	118.32	110.41
1	A	89	LYS	N-CA-C	-6.35	105.11	112.86
1	A	255	ILE	N-CA-C	6.31	117.06	110.62
1	B	48	ILE	N-CA-C	6.30	116.47	110.42
1	B	297	GLU	N-CA-C	6.27	119.32	110.23
1	A	313	ASP	N-CA-C	-6.15	104.84	113.20
1	B	503	TYR	N-CA-C	6.13	120.79	113.12
1	B	283	ASP	N-CA-C	-5.99	105.06	112.90
1	B	279	THR	N-CA-C	-5.98	105.85	113.02
1	A	150	VAL	N-CA-C	5.97	116.16	110.42
1	A	29	SER	N-CA-C	-5.97	107.23	114.75
1	A	382	HIS	N-CA-C	5.94	119.84	111.52
1	B	332	SER	N-CA-C	5.74	119.27	110.20
1	B	158	CYS	N-CA-C	5.67	122.88	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	258	ASP	N-CA-C	5.61	117.47	111.36
1	A	136	ARG	N-CA-C	5.48	117.83	108.90
1	B	175	ARG	N-CA-C	-5.47	106.53	113.43
1	B	28	GLN	N-CA-C	5.40	118.53	109.95
1	A	314	GLY	N-CA-C	5.37	121.38	112.89
1	B	495	LEU	N-CA-C	-5.34	105.56	111.71
1	B	336	THR	N-CA-C	-5.14	105.73	112.41
1	A	421	ASP	N-CA-C	5.13	118.09	110.14
1	A	308	TYR	N-CA-C	5.09	118.02	110.48
1	B	280	ASN	N-CA-C	5.06	117.56	109.96
1	B	328	VAL	CA-C-O	5.01	121.93	119.12

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	3875	0	3823	242	0
1	B	3879	0	3827	249	0
2	A	24	0	32	9	0
2	B	12	0	16	5	0
3	A	16	0	11	0	0
3	B	16	0	11	1	0
4	A	43	0	0	2	0
4	B	38	0	0	1	0
All	All	7903	0	7720	476	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 31.

All (476) 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:212:MET:HE2	1:A:450:MET:HE1	1.22	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:480:ILE:HD13	1:B:491:GLU:HG3	1.29	1.15
1:A:439:ARG:HA	1:B:411:MET:HE1	1.33	1.09
1:A:238:LYS:HE2	1:A:364:GLU:HG2	1.32	1.08
1:B:182:ARG:HH11	1:B:182:ARG:HG2	1.19	1.07
1:B:125:MET:HE2	1:B:136:ARG:HG2	1.33	1.06
1:B:330:ARG:H	1:B:330:ARG:HD3	1.15	1.05
1:B:330:ARG:H	1:B:330:ARG:CD	1.63	1.02
1:B:355:ILE:HG21	1:B:496:MET:HE3	1.42	1.01
1:A:284:VAL:HG21	1:A:450:MET:HE3	1.48	0.93
1:A:260:MET:HE1	1:A:340:VAL:HG21	1.53	0.90
1:B:171:LYS:HB2	1:B:172:PRO:HA	1.51	0.89
1:A:146:VAL:HG12	1:A:150:VAL:HG23	1.55	0.89
1:A:260:MET:HE2	1:A:465:PHE:CE1	2.08	0.89
1:B:137:LEU:HD21	1:B:168:ILE:HD11	1.52	0.89
1:B:404:MET:HE3	1:B:420:LEU:CD1	2.02	0.89
1:A:171:LYS:HB2	1:A:172:PRO:HA	1.55	0.88
1:B:243:THR:HG21	1:B:247:GLY:CA	2.04	0.88
1:A:138:PHE:HB2	1:A:167:ILE:HD12	1.56	0.88
1:A:442:LEU:HD22	1:B:411:MET:HE2	1.55	0.87
1:B:125:MET:CE	1:B:136:ARG:HG2	2.02	0.87
1:A:355:ILE:HG21	1:A:496:MET:HE3	1.57	0.86
1:A:91:THR:HB	1:A:92:PRO:HD2	1.55	0.86
1:A:411:MET:HE1	1:B:439:ARG:HA	1.55	0.85
1:B:355:ILE:HG21	1:B:496:MET:CE	2.07	0.85
1:B:330:ARG:HD3	1:B:330:ARG:N	1.92	0.84
1:B:138:PHE:HB2	1:B:167:ILE:HD12	1.59	0.84
1:B:243:THR:HG21	1:B:247:GLY:HA2	1.60	0.84
1:A:439:ARG:HA	1:B:411:MET:CE	2.08	0.82
1:A:27:VAL:N	1:A:30:ASP:HB2	1.95	0.81
1:B:480:ILE:CD1	1:B:491:GLU:HG3	2.08	0.81
1:A:260:MET:HE1	1:A:340:VAL:CG2	2.10	0.80
1:B:407:LYS:HE2	1:B:412:PHE:O	1.82	0.80
1:A:407:LYS:HE2	1:A:412:PHE:O	1.82	0.80
1:B:303:VAL:HG22	1:B:340:VAL:HG22	1.64	0.80
1:B:146:VAL:HG12	1:B:150:VAL:HG23	1.63	0.79
1:A:459:ARG:HH11	1:A:459:ARG:CG	1.95	0.79
1:A:292:LEU:HA	1:A:295:ILE:HD12	1.64	0.78
1:B:238:LYS:HE2	1:B:364:GLU:HG2	1.64	0.78
1:B:287:GLU:O	1:B:291:VAL:HG23	1.84	0.78
1:A:27:VAL:HG12	1:A:28:GLN:HG3	1.65	0.77
1:B:146:VAL:HG12	1:B:150:VAL:CG2	2.15	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:ARG:HG2	1:A:463:ARG:HH11	1.49	0.76
1:B:320:LYS:HE2	1:B:324:ASP:O	1.87	0.75
1:A:212:MET:CE	1:A:450:MET:HE1	2.11	0.74
1:B:341:VAL:HG21	1:B:499:VAL:HG21	1.69	0.74
1:A:176:ASP:H	1:A:179:SER:HB2	1.51	0.74
1:A:264:LEU:HD12	1:A:267:MET:HE2	1.69	0.74
1:B:182:ARG:HH11	1:B:182:ARG:CG	2.00	0.74
1:A:71:ALA:HB3	1:A:107:TYR:OH	1.88	0.74
1:A:411:MET:CE	1:B:439:ARG:HA	2.18	0.74
1:A:264:LEU:HD12	1:A:267:MET:CE	2.18	0.73
1:B:330:ARG:CD	1:B:330:ARG:N	2.47	0.73
1:A:459:ARG:HH11	1:A:459:ARG:HG3	1.51	0.73
1:B:480:ILE:HD13	1:B:491:GLU:CG	2.13	0.73
1:A:138:PHE:HB2	1:A:167:ILE:CD1	2.19	0.73
1:A:56:PHE:HA	1:A:61:LEU:HD12	1.71	0.73
1:A:56:PHE:CE2	1:A:101:PHE:HB2	2.24	0.72
1:B:346:ASN:O	1:B:350:ASP:HB2	1.90	0.72
1:B:211:LEU:CD2	1:B:392:ILE:HD13	2.20	0.71
1:B:273:MET:HA	1:B:291:VAL:CG2	2.19	0.71
1:B:325:ASP:HB3	1:B:328:VAL:HG23	1.70	0.71
1:A:273:MET:HE2	1:A:276:PRO:HD3	1.71	0.71
1:B:325:ASP:HB3	1:B:328:VAL:CG2	2.20	0.71
1:B:281:SER:HA	1:B:450:MET:HB2	1.73	0.71
1:B:459:ARG:HH11	1:B:459:ARG:HG2	1.54	0.71
1:A:60:LEU:HB3	1:A:442:LEU:HD11	1.73	0.71
1:B:405:MET:HE3	1:B:416:GLU:O	1.91	0.71
1:B:71:ALA:HB3	1:B:107:TYR:OH	1.90	0.70
1:B:169:VAL:HG12	1:B:170:GLU:N	2.07	0.70
1:A:260:MET:HE2	1:A:465:PHE:HE1	1.51	0.70
1:A:240:PRO:HD3	1:A:365:ARG:HB2	1.73	0.70
1:B:480:ILE:HG21	1:B:489:PRO:HB3	1.74	0.69
1:B:459:ARG:HG2	1:B:459:ARG:NH1	2.08	0.68
1:B:273:MET:HA	1:B:291:VAL:HG21	1.75	0.68
1:A:48:ILE:N	1:A:48:ILE:HD13	2.08	0.68
1:A:411:MET:HE1	1:B:439:ARG:CA	2.23	0.68
1:A:273:MET:HE3	1:A:287:GLU:CB	2.24	0.68
1:A:430:ASN:OD1	1:A:430:ASN:N	2.26	0.68
1:B:243:THR:HG21	1:B:247:GLY:N	2.08	0.68
1:B:469:LEU:HA	1:B:472:ILE:HD12	1.75	0.68
1:A:393:ARG:NH1	1:A:397:ASN:O	2.27	0.67
1:A:456:ASP:N	1:A:456:ASP:OD1	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:ARG:HG2	1:A:463:ARG:NH1	2.08	0.67
1:B:256:ILE:HA	1:B:260:MET:HB2	1.75	0.67
1:A:56:PHE:CA	1:A:61:LEU:HD12	2.25	0.67
1:A:146:VAL:HG12	1:A:150:VAL:CG2	2.23	0.67
1:A:273:MET:HA	1:A:291:VAL:HG21	1.76	0.67
1:A:300:ALA:HB1	2:A:1507:GOL:H2	1.77	0.67
1:B:137:LEU:HD21	1:B:168:ILE:CD1	2.24	0.67
1:B:144:PRO:HG2	1:B:172:PRO:HD2	1.74	0.67
1:A:393:ARG:HD2	1:A:396:PRO:O	1.95	0.66
1:B:95:LYS:NZ	1:B:99:GLU:OE2	2.28	0.66
1:A:442:LEU:HD22	1:B:411:MET:CE	2.24	0.66
1:B:125:MET:O	1:B:131:GLY:HA3	1.96	0.66
1:B:177:LEU:HB2	1:B:462:TRP:HB3	1.77	0.66
1:A:81:ARG:O	1:A:85:GLU:HB2	1.95	0.66
1:A:53:TRP:O	1:A:56:PHE:HB3	1.97	0.65
1:A:92:PRO:O	1:A:95:LYS:HG3	1.96	0.65
1:A:473:GLU:O	1:A:476:LYS:HE2	1.96	0.65
1:B:333:THR:OG1	1:B:476:LYS:HD3	1.95	0.65
1:A:154:ILE:O	1:A:159:MET:HG2	1.95	0.65
1:B:278:SER:OG	1:B:280:ASN:HB2	1.95	0.65
1:A:435:ASP:HB3	1:A:438:GLU:HG3	1.79	0.64
1:A:342:LEU:HD11	1:A:356:LEU:CD1	2.27	0.64
1:A:238:LYS:HE2	1:A:364:GLU:CG	2.19	0.64
1:A:309:VAL:CG2	1:A:483:ILE:HD13	2.27	0.64
1:B:144:PRO:CG	1:B:172:PRO:HD2	2.27	0.64
1:A:312:PRO:C	1:A:314:GLY:H	2.04	0.64
1:A:231:ALA:HB2	1:A:374:HIS:CD2	2.33	0.64
1:B:55:LEU:HD23	1:B:60:LEU:HD12	1.79	0.64
1:B:450:MET:HE3	1:B:451:HIS:CE1	2.32	0.64
1:A:198:ARG:HG3	1:A:198:ARG:HH11	1.63	0.63
1:B:404:MET:HE3	1:B:420:LEU:HD12	1.80	0.63
1:A:213:VAL:O	1:A:217:ALA:HB3	1.99	0.63
1:B:255:ILE:HG12	1:B:358:CYS:SG	2.39	0.63
1:B:355:ILE:CG2	1:B:496:MET:HE3	2.25	0.63
1:A:271:VAL:HG21	1:A:369:VAL:HG11	1.81	0.62
1:A:27:VAL:HG12	1:A:28:GLN:H	1.64	0.62
1:A:408:LYS:HD3	1:A:412:PHE:CG	2.34	0.62
1:A:466:THR:HB	1:A:467:PRO:HD3	1.82	0.62
1:A:346:ASN:O	1:A:350:ASP:HB2	2.00	0.62
1:B:211:LEU:HD23	1:B:392:ILE:HD13	1.80	0.62
1:B:323:LEU:HD23	1:B:328:VAL:HG11	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:MET:HE3	1:A:287:GLU:HB3	1.81	0.62
1:B:166:ARG:NH1	1:B:444:VAL:O	2.30	0.62
1:B:183:LEU:HG	1:B:187:ILE:HD11	1.82	0.62
1:A:125:MET:O	1:A:131:GLY:HA3	2.00	0.62
1:A:146:VAL:CG1	1:A:150:VAL:HG23	2.29	0.61
1:A:166:ARG:NH1	1:A:444:VAL:O	2.34	0.61
1:A:411:MET:HE3	1:B:442:LEU:HD23	1.82	0.61
1:B:169:VAL:HG12	1:B:170:GLU:H	1.64	0.61
1:A:305:LEU:HD11	2:A:1507:GOL:H12	1.83	0.60
1:A:342:LEU:HD11	1:A:356:LEU:HD11	1.82	0.60
1:B:138:PHE:HB2	1:B:167:ILE:CD1	2.30	0.60
1:A:49:TYR:O	1:A:52:ILE:HG22	2.02	0.60
1:A:468:LEU:O	1:A:472:ILE:HG13	2.01	0.60
1:B:129:HIS:O	1:B:130:LEU:HB2	2.01	0.60
1:A:307:GLN:HA	1:A:336:THR:OG1	2.02	0.60
1:B:171:LYS:CB	1:B:172:PRO:HA	2.18	0.60
1:A:287:GLU:O	1:A:291:VAL:HG23	2.02	0.60
1:A:381:PHE:O	1:A:384:GLN:HG3	2.02	0.60
1:B:35:ILE:HD11	1:B:125:MET:CE	2.32	0.60
1:B:381:PHE:O	1:B:384:GLN:HG3	2.01	0.60
1:B:405:MET:HE3	1:B:405:MET:HA	1.82	0.60
1:B:273:MET:CA	1:B:291:VAL:HG21	2.31	0.59
1:A:122:ASN:O	1:A:125:MET:N	2.36	0.59
1:B:44:ALA:HA	1:B:48:ILE:HB	1.85	0.59
1:A:99:GLU:OE2	1:A:99:GLU:HA	2.02	0.58
1:A:268:LEU:HD23	1:A:268:LEU:C	2.29	0.58
1:A:73:SER:O	1:A:75:LEU:HG	2.03	0.58
1:A:144:PRO:HG3	1:A:172:PRO:HD2	1.86	0.58
1:B:307:GLN:O	1:B:481:PRO:HA	2.04	0.58
1:A:170:GLU:HG2	4:A:2007:HOH:O	2.04	0.57
1:A:247:GLY:O	1:A:328:VAL:HG22	2.05	0.57
1:B:37:MET:HE1	1:B:70:TYR:CD1	2.39	0.57
1:B:202:TYR:HA	1:B:205:LYS:HD2	1.86	0.57
1:B:325:ASP:O	1:B:328:VAL:HB	2.04	0.57
1:A:273:MET:HA	1:A:291:VAL:CG2	2.35	0.57
1:B:243:THR:CG2	1:B:247:GLY:N	2.67	0.57
1:B:197:TYR:CD2	1:B:444:VAL:HG22	2.38	0.57
1:B:235:LEU:HD12	1:B:356:LEU:HD23	1.86	0.57
1:B:456:ASP:OD1	1:B:456:ASP:N	2.36	0.57
1:B:492:ALA:O	1:B:496:MET:HG3	2.05	0.57
1:A:260:MET:HE1	1:A:340:VAL:CB	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:GLY:HA3	4:B:2002:HOH:O	2.04	0.56
1:A:171:LYS:CB	1:A:172:PRO:HA	2.24	0.56
1:B:114:ASP:O	1:B:118:TYR:HD1	1.87	0.56
1:B:182:ARG:HG2	1:B:182:ARG:NH1	1.99	0.56
1:B:190:LEU:HB2	1:B:191:PHE:CD1	2.40	0.56
1:A:35:ILE:HD12	1:A:35:ILE:H	1.70	0.56
1:B:183:LEU:HG	1:B:187:ILE:CD1	2.36	0.56
1:A:240:PRO:HD3	1:A:365:ARG:CB	2.35	0.56
1:A:129:HIS:O	1:A:130:LEU:HB2	2.06	0.55
1:A:298:VAL:HA	2:A:1506:GOL:O1	2.06	0.55
1:B:426:ASN:O	1:B:429:LYS:HG2	2.07	0.55
1:A:88:PHE:O	1:A:89:LYS:HB3	2.07	0.55
1:A:264:LEU:CD2	1:A:358:CYS:HB3	2.37	0.55
1:A:132:SER:O	1:A:162:ILE:HD12	2.06	0.55
1:A:305:LEU:CD1	2:A:1507:GOL:H12	2.37	0.55
1:A:342:LEU:CD1	1:A:356:LEU:HD11	2.36	0.55
1:B:49:TYR:CB	1:B:50:PRO:HD3	2.36	0.55
1:A:307:GLN:O	1:A:481:PRO:HA	2.06	0.55
1:B:405:MET:CE	1:B:416:GLU:C	2.80	0.55
1:A:434:PRO:HD2	1:A:439:ARG:NH2	2.22	0.55
1:A:273:MET:CE	1:A:276:PRO:HD3	2.37	0.54
1:A:282:ASP:OD1	1:A:285:ARG:NH1	2.29	0.54
1:B:177:LEU:HD12	1:B:177:LEU:O	2.07	0.54
1:B:262:ASN:OD1	1:B:262:ASN:N	2.35	0.54
1:B:235:LEU:HB2	1:B:356:LEU:HD23	1.89	0.54
1:B:273:MET:CE	1:B:276:PRO:HD3	2.38	0.54
1:B:206:GLU:OE1	1:B:439:ARG:NH2	2.39	0.54
1:B:264:LEU:CD2	1:B:358:CYS:HB3	2.36	0.54
1:B:468:LEU:O	1:B:472:ILE:HD12	2.07	0.54
1:A:198:ARG:HG3	1:A:198:ARG:NH1	2.23	0.54
1:B:298:VAL:HG22	1:B:342:LEU:CD2	2.38	0.54
1:A:282:ASP:OD1	1:A:285:ARG:HD3	2.08	0.53
1:A:62:PRO:HD2	1:A:65:THR:OG1	2.07	0.53
1:A:121:LEU:O	1:A:124:HIS:HB3	2.09	0.53
1:B:34:PHE:CE1	1:B:137:LEU:HD22	2.43	0.53
1:B:298:VAL:HG22	1:B:342:LEU:HD21	1.89	0.53
1:A:133:GLN:HA	1:A:162:ILE:HD12	1.90	0.53
1:A:211:LEU:HD23	1:A:392:ILE:HD13	1.91	0.53
1:B:225:TRP:CZ2	1:B:371:LEU:HD11	2.43	0.53
1:A:138:PHE:CD1	1:A:167:ILE:CD1	2.92	0.53
1:B:432:LYS:NZ	2:B:1506:GOL:H11	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:PHE:HB3	1:B:275:LYS:HE3	1.90	0.52
1:A:27:VAL:HG12	1:A:28:GLN:CG	2.38	0.52
1:A:308:TYR:HB3	1:A:482:TYR:CZ	2.45	0.52
1:B:95:LYS:O	1:B:99:GLU:HG2	2.09	0.52
1:B:211:LEU:HD21	1:B:392:ILE:HD13	1.89	0.52
1:B:325:ASP:OD1	1:B:326:PRO:HD2	2.10	0.52
1:A:367:ALA:HB3	1:A:395:GLN:HG2	1.91	0.52
1:A:85:GLU:N	1:A:86:PRO:CD	2.73	0.52
1:A:132:SER:C	1:A:162:ILE:HD12	2.35	0.52
1:B:239:GLU:HG2	1:B:241:PHE:CZ	2.44	0.52
1:B:309:VAL:CG2	1:B:483:ILE:HD13	2.39	0.52
1:B:260:MET:HG3	1:B:358:CYS:SG	2.50	0.52
1:A:117:SER:HA	1:A:120:ARG:CZ	2.40	0.52
1:A:117:SER:O	1:A:120:ARG:HB2	2.10	0.52
1:B:242:GLY:O	1:B:244:GLU:HG3	2.10	0.52
1:B:260:MET:CE	1:B:465:PHE:CE1	2.93	0.52
1:A:137:LEU:CD1	1:A:164:TRP:CZ3	2.93	0.51
1:A:375:ASP:OD1	1:A:375:ASP:N	2.43	0.51
1:B:269:CYS:SG	1:B:292:LEU:HD21	2.50	0.51
1:B:389:GLU:OE1	1:B:403:LYS:HE2	2.11	0.51
1:B:84:SER:C	1:B:86:PRO:HD2	2.35	0.51
1:B:170:GLU:HG3	1:B:199:ILE:HB	1.92	0.51
1:A:271:VAL:CG2	1:A:369:VAL:HG11	2.40	0.51
1:B:468:LEU:HG	1:B:472:ILE:HD11	1.92	0.51
1:A:480:ILE:HG21	1:A:489:PRO:HB3	1.92	0.51
1:B:49:TYR:HB2	1:B:50:PRO:HD3	1.93	0.51
1:A:309:VAL:HG23	1:A:483:ILE:HD13	1.92	0.51
1:A:206:GLU:OE2	1:B:407:LYS:NZ	2.42	0.51
1:A:207:MET:HB2	1:A:394:VAL:HA	1.91	0.51
1:A:175:ARG:NH2	1:A:252:GLU:O	2.33	0.51
1:B:169:VAL:CG1	1:B:170:GLU:N	2.74	0.51
1:A:110:GLY:HA3	1:A:117:SER:HB3	1.93	0.50
1:A:42:ASP:O	1:A:46:LYS:HB2	2.10	0.50
1:B:412:PHE:N	1:B:412:PHE:CD2	2.79	0.50
1:A:108:VAL:HG11	1:A:121:LEU:HA	1.92	0.50
1:A:300:ALA:HB1	2:A:1507:GOL:C2	2.42	0.50
1:A:49:TYR:CB	1:A:50:PRO:HD3	2.42	0.50
1:A:235:LEU:CD1	1:A:268:LEU:HB2	2.42	0.50
1:A:260:MET:CE	1:A:465:PHE:HE1	2.22	0.50
1:A:264:LEU:CD1	1:A:267:MET:HE2	2.41	0.50
1:A:302:ASN:OD1	1:A:498:ARG:NH2	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:ILE:HD12	1:B:35:ILE:H	1.76	0.50
1:B:323:LEU:HD23	1:B:328:VAL:CG1	2.42	0.50
1:A:144:PRO:CG	1:A:172:PRO:HD2	2.41	0.50
1:A:393:ARG:CD	1:A:396:PRO:O	2.60	0.50
1:A:372:GLN:HA	1:A:389:GLU:HG2	1.93	0.50
1:B:304:VAL:HB	1:B:339:ALA:HB3	1.93	0.50
1:A:133:GLN:HA	1:A:162:ILE:HB	1.93	0.50
1:A:300:ALA:CB	2:A:1507:GOL:H2	2.42	0.50
1:A:440:LEU:O	1:A:443:ASP:HB2	2.12	0.50
1:B:57:ARG:HA	1:B:97:LYS:HD3	1.93	0.49
1:B:333:THR:HB	1:B:477:PRO:O	2.12	0.49
1:B:438:GLU:O	1:B:442:LEU:HB2	2.12	0.49
1:A:231:ALA:HB2	1:A:374:HIS:HD2	1.75	0.49
1:A:393:ARG:NH1	1:A:397:ASN:HB3	2.28	0.49
1:B:466:THR:HB	1:B:467:PRO:HD3	1.95	0.49
1:A:190:LEU:HB2	1:A:191:PHE:CD1	2.48	0.49
1:A:241:PHE:O	1:A:361:ALA:N	2.45	0.49
1:A:404:MET:HE2	1:A:405:MET:O	2.12	0.49
1:B:181:ASP:OD2	1:B:459:ARG:NH1	2.46	0.49
1:B:381:PHE:CD1	1:B:415:PRO:HG3	2.48	0.49
1:B:37:MET:CE	1:B:70:TYR:CD1	2.95	0.49
1:B:250:PHE:O	1:B:253:PHE:N	2.39	0.49
1:A:91:THR:CB	1:A:92:PRO:HD2	2.33	0.49
1:A:492:ALA:O	1:A:496:MET:HG3	2.13	0.49
1:B:197:TYR:OH	1:B:447:GLY:HA2	2.12	0.49
1:B:260:MET:CE	1:B:465:PHE:CZ	2.95	0.49
1:B:273:MET:HA	1:B:291:VAL:HG22	1.94	0.49
1:B:476:LYS:N	1:B:477:PRO:HD3	2.27	0.49
1:A:312:PRO:C	1:A:314:GLY:N	2.70	0.49
1:B:393:ARG:HD2	1:B:396:PRO:O	2.13	0.49
1:A:311:ASN:O	1:A:314:GLY:HA3	2.13	0.49
1:A:137:LEU:HD22	1:A:445:PHE:CZ	2.48	0.48
1:A:150:VAL:O	1:A:154:ILE:HG13	2.12	0.48
1:A:247:GLY:O	1:A:251:ASP:HB2	2.13	0.48
1:A:411:MET:CE	1:B:442:LEU:HD23	2.42	0.48
1:A:375:ASP:O	1:B:219:ARG:NH1	2.47	0.48
1:A:260:MET:HE2	1:A:465:PHE:CZ	2.48	0.48
1:B:237:PHE:CG	1:B:259:VAL:HG11	2.48	0.48
1:A:84:SER:C	1:A:86:PRO:HD2	2.38	0.48
1:A:235:LEU:HD11	1:A:268:LEU:HB2	1.95	0.48
1:B:454:ARG:O	1:B:457:GLU:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:LEU:HD21	1:A:358:CYS:HB3	1.95	0.48
1:A:273:MET:HE3	1:A:287:GLU:HB2	1.96	0.48
1:A:459:ARG:HG3	1:A:459:ARG:NH1	2.25	0.48
1:A:168:ILE:HD11	1:A:444:VAL:HG21	1.95	0.48
1:A:477:PRO:HG3	2:A:1507:GOL:C3	2.43	0.48
1:A:27:VAL:HG12	1:A:28:GLN:N	2.28	0.47
1:A:342:LEU:HD12	1:A:356:LEU:HG	1.96	0.47
1:B:42:ASP:O	1:B:46:LYS:HG3	2.14	0.47
1:B:235:LEU:CD1	1:B:356:LEU:HD23	2.44	0.47
1:A:271:VAL:HG21	1:A:369:VAL:CG1	2.44	0.47
1:B:306:GLY:HA2	1:B:480:ILE:O	2.13	0.47
1:B:400:VAL:HG13	1:B:422:LEU:HB3	1.95	0.47
1:A:342:LEU:CD1	1:A:356:LEU:CD1	2.92	0.47
1:B:333:THR:CG2	1:B:476:LYS:HD3	2.44	0.47
1:A:27:VAL:CG1	1:A:28:GLN:H	2.24	0.47
1:A:49:TYR:HB2	1:A:50:PRO:HD3	1.96	0.47
1:A:183:LEU:O	1:A:187:ILE:HG13	2.14	0.47
1:B:85:GLU:N	1:B:86:PRO:CD	2.77	0.47
1:B:298:VAL:HG23	1:B:464:ILE:HG22	1.95	0.47
1:B:137:LEU:CD2	1:B:168:ILE:CD1	2.91	0.47
1:B:235:LEU:HD12	1:B:356:LEU:CD2	2.45	0.47
1:B:302:ASN:O	1:B:495:LEU:HD21	2.14	0.47
1:B:353:PRO:CB	1:B:499:VAL:CG1	2.93	0.47
1:A:137:LEU:CD1	1:A:164:TRP:HZ3	2.26	0.47
1:A:138:PHE:CD1	1:A:167:ILE:HD11	2.49	0.47
1:A:237:PHE:CD1	1:A:237:PHE:C	2.93	0.47
1:B:169:VAL:CG1	1:B:170:GLU:H	2.27	0.47
1:B:268:LEU:C	1:B:268:LEU:HD23	2.39	0.47
1:B:408:LYS:HE2	1:B:412:PHE:CD1	2.50	0.47
1:A:409:PRO:HA	1:B:206:GLU:OE2	2.15	0.47
1:B:175:ARG:O	1:B:462:TRP:CZ3	2.67	0.47
1:B:495:LEU:O	1:B:499:VAL:HG23	2.14	0.47
1:A:260:MET:HA	1:A:264:LEU:HB2	1.97	0.47
1:A:463:ARG:HH11	1:A:463:ARG:CG	2.20	0.47
1:B:166:ARG:C	1:B:167:ILE:HD13	2.40	0.47
1:B:360:LYS:NZ	3:B:1508:BG6:O3	2.48	0.47
1:A:70:TYR:HD2	1:A:121:LEU:HD13	1.80	0.47
1:B:35:ILE:CG1	1:B:125:MET:HE3	2.45	0.47
1:B:182:ARG:CG	1:B:182:ARG:NH1	2.66	0.46
1:A:123:SER:HA	1:A:126:ASN:HB2	1.97	0.46
1:B:289:VAL:O	1:B:293:LYS:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ASP:OD2	1:A:179:SER:HB2	2.15	0.46
1:A:260:MET:CE	1:A:340:VAL:HB	2.46	0.46
1:A:298:VAL:HG12	1:A:299:GLN:N	2.30	0.46
1:A:408:LYS:HD2	1:A:414:ASN:HD21	1.80	0.46
1:B:273:MET:HE2	1:B:276:PRO:HD3	1.97	0.46
1:B:304:VAL:HG23	1:B:495:LEU:HD22	1.96	0.46
1:B:405:MET:HE3	1:B:416:GLU:C	2.40	0.46
1:A:238:LYS:CE	1:A:364:GLU:HG2	2.24	0.46
1:A:355:ILE:HD13	1:A:496:MET:CE	2.45	0.46
1:A:398:GLU:CD	1:A:398:GLU:H	2.23	0.46
1:B:46:LYS:HD3	1:B:87:PHE:CE2	2.51	0.46
1:B:227:ARG:HA	1:B:352:VAL:HG21	1.97	0.46
1:B:432:LYS:HZ2	2:B:1506:GOL:H11	1.80	0.46
1:A:260:MET:HE1	1:A:340:VAL:HB	1.98	0.46
1:A:295:ILE:HG12	1:A:344:VAL:HG22	1.98	0.46
1:B:35:ILE:HD12	1:B:35:ILE:N	2.31	0.46
1:A:81:ARG:HH12	1:A:99:GLU:CD	2.23	0.46
1:B:209:GLN:NE2	1:B:439:ARG:HD2	2.31	0.46
1:A:235:LEU:CD2	1:A:369:VAL:HG13	2.46	0.45
1:B:35:ILE:HD12	1:B:137:LEU:O	2.16	0.45
1:A:72:ARG:HH11	1:A:72:ARG:CG	2.30	0.45
1:A:80:ILE:HG22	1:A:81:ARG:N	2.32	0.45
1:A:405:MET:HE3	1:A:415:PRO:HB2	1.98	0.45
1:A:231:ALA:CB	1:A:374:HIS:CD2	3.00	0.45
1:B:144:PRO:HA	1:B:147:TYR:CG	2.52	0.45
1:B:337:PHE:CE2	1:B:489:PRO:HD2	2.52	0.45
1:B:400:VAL:HG13	1:B:400:VAL:O	2.16	0.45
1:A:220:ILE:HB	1:B:388:ASN:OD1	2.16	0.45
1:B:238:LYS:HG2	1:B:362:LEU:HD12	1.99	0.45
1:B:292:LEU:O	1:B:463:ARG:NH2	2.49	0.45
1:A:269:CYS:O	1:A:273:MET:HB3	2.17	0.45
1:A:477:PRO:HG3	2:A:1507:GOL:H32	1.99	0.45
1:A:385:CYS:HA	1:A:405:MET:HE2	1.99	0.45
1:B:57:ARG:NH2	2:B:1506:GOL:O1	2.49	0.45
1:A:109:ALA:O	1:A:120:ARG:NH1	2.50	0.44
1:A:370:ARG:C	1:A:371:LEU:HD23	2.41	0.44
1:A:408:LYS:HB3	1:A:412:PHE:HB2	1.98	0.44
1:B:436:ALA:O	1:B:440:LEU:HB2	2.17	0.44
1:B:260:MET:HE1	1:B:465:PHE:CE1	2.51	0.44
1:B:423:THR:HG22	1:B:425:GLY:H	1.81	0.44
1:B:432:LYS:NZ	2:B:1506:GOL:C1	2.79	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ARG:HB2	1:A:453:VAL:O	2.18	0.44
1:A:297:GLU:O	2:A:1506:GOL:H2	2.17	0.44
1:A:355:ILE:C	1:A:356:LEU:HD23	2.42	0.44
1:B:251:ASP:C	1:B:253:PHE:H	2.25	0.44
1:A:125:MET:HA	1:A:128:LEU:HG	1.99	0.44
1:A:230:ILE:HG12	1:A:373:PHE:CE1	2.53	0.44
1:B:176:ASP:HB2	1:B:179:SER:OG	2.18	0.44
1:B:315:GLU:O	1:B:318:ALA:HB3	2.17	0.44
1:B:450:MET:HE2	1:B:450:MET:HB3	1.70	0.44
1:A:55:LEU:HD22	1:A:442:LEU:CD1	2.48	0.44
1:A:306:GLY:HA2	1:A:480:ILE:O	2.18	0.44
1:B:36:ILE:CD1	1:B:48:ILE:HG22	2.48	0.43
1:B:172:PRO:HG3	1:B:253:PHE:CZ	2.53	0.43
1:B:299:GLN:C	1:B:301:ASN:H	2.26	0.43
1:B:305:LEU:HB3	1:B:335:ALA:HB2	1.98	0.43
1:A:273:MET:CA	1:A:291:VAL:HG21	2.45	0.43
1:A:480:ILE:CG2	1:A:489:PRO:HB3	2.48	0.43
1:B:261:GLN:HE21	1:B:261:GLN:HB3	1.66	0.43
1:B:380:ILE:HG12	1:B:381:PHE:CE2	2.54	0.43
1:B:432:LYS:HZ2	2:B:1506:GOL:C1	2.32	0.43
1:B:142:LEU:HB3	1:B:143:PRO:HD2	2.00	0.43
1:A:88:PHE:HD2	1:A:88:PHE:N	2.17	0.43
1:B:36:ILE:HD13	1:B:48:ILE:HG22	1.99	0.43
1:B:172:PRO:HG3	1:B:253:PHE:HZ	1.84	0.43
1:B:235:LEU:CD1	1:B:356:LEU:CD2	2.97	0.43
1:A:30:ASP:O	1:A:32:HIS:HD2	2.01	0.43
1:B:53:TRP:CD1	1:B:88:PHE:HB3	2.54	0.43
1:A:281:SER:OG	1:A:450:MET:N	2.52	0.43
1:A:308:TYR:HB3	1:A:482:TYR:CE2	2.53	0.42
1:B:35:ILE:HD11	1:B:125:MET:HE1	1.99	0.42
1:A:146:VAL:HG12	1:A:146:VAL:O	2.18	0.42
1:B:309:VAL:HG23	1:B:483:ILE:HD13	2.01	0.42
1:A:34:PHE:CE1	1:A:137:LEU:HD23	2.53	0.42
1:A:431:VAL:HG12	1:A:432:LYS:N	2.35	0.42
1:B:35:ILE:HG12	1:B:125:MET:CE	2.49	0.42
1:A:139:TYR:CE1	1:A:168:ILE:HG21	2.53	0.42
1:A:212:MET:HE2	1:A:450:MET:CE	2.16	0.42
1:B:150:VAL:HG12	1:B:154:ILE:HD12	2.00	0.42
1:B:459:ARG:HH11	1:B:459:ARG:CG	2.23	0.42
1:B:303:VAL:CG1	1:B:304:VAL:N	2.82	0.42
1:A:487:ARG:NH2	1:A:493:ASP:OD1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:VAL:O	1:B:304:VAL:HG12	2.19	0.42
1:A:91:THR:HB	1:A:92:PRO:CD	2.38	0.42
1:A:199:ILE:HG22	4:A:2007:HOH:O	2.19	0.42
1:A:404:MET:HE1	1:A:406:THR:HB	2.01	0.42
1:A:407:LYS:NZ	1:B:206:GLU:OE2	2.53	0.42
1:B:35:ILE:CD1	1:B:125:MET:CE	2.98	0.42
1:B:298:VAL:HG23	1:B:464:ILE:CG2	2.50	0.42
1:A:34:PHE:CD1	1:A:137:LEU:HD23	2.55	0.42
1:A:278:SER:C	1:A:280:ASN:H	2.28	0.42
1:B:159:MET:HE3	1:B:159:MET:HB3	1.75	0.42
1:B:405:MET:HE1	1:B:416:GLU:C	2.44	0.42
1:B:493:ASP:O	1:B:496:MET:HB2	2.19	0.42
1:B:307:GLN:HB2	1:B:479:PRO:HB2	2.01	0.41
1:B:325:ASP:HA	1:B:326:PRO:HD3	1.89	0.41
1:A:155:HIS:CD2	1:A:190:LEU:HB3	2.55	0.41
1:B:35:ILE:HD11	1:B:125:MET:HE3	2.00	0.41
1:B:49:TYR:CB	1:B:50:PRO:CD	2.98	0.41
1:B:250:PHE:O	1:B:253:PHE:HB2	2.19	0.41
1:B:260:MET:HE2	1:B:260:MET:HB3	1.74	0.41
1:B:466:THR:CB	1:B:467:PRO:HD3	2.51	0.41
1:A:72:ARG:HH11	1:A:72:ARG:HG2	1.85	0.41
1:A:85:GLU:N	1:A:86:PRO:HD2	2.36	0.41
1:A:325:ASP:HA	1:A:326:PRO:HD3	1.80	0.41
1:B:55:LEU:CD2	1:B:60:LEU:HD12	2.48	0.41
1:B:265:LEU:O	1:B:268:LEU:HB3	2.20	0.41
1:B:425:GLY:O	1:B:429:LYS:HD3	2.20	0.41
1:B:435:ASP:OD2	1:B:437:TYR:HB2	2.20	0.41
1:B:480:ILE:CG2	1:B:489:PRO:HB3	2.46	0.41
1:B:31:THR:HA	1:B:64:ASN:HB2	2.03	0.41
1:B:33:ILE:HD13	1:B:125:MET:HG2	2.02	0.41
1:B:171:LYS:HA	1:B:172:PRO:C	2.44	0.41
1:B:278:SER:C	1:B:280:ASN:H	2.26	0.41
1:A:137:LEU:HD13	1:A:164:TRP:CZ3	2.54	0.41
1:A:325:ASP:OD2	1:A:327:THR:HG23	2.19	0.41
1:B:408:LYS:HG2	1:B:412:PHE:CG	2.55	0.41
1:A:33:ILE:HD13	1:A:125:MET:HB3	2.02	0.41
1:A:238:LYS:HB3	1:A:238:LYS:HE3	1.95	0.41
1:B:36:ILE:HD13	1:B:48:ILE:CG2	2.51	0.41
1:B:400:VAL:CG1	1:B:422:LEU:HB3	2.51	0.41
1:B:466:THR:CB	1:B:467:PRO:CD	2.99	0.41
1:A:57:ARG:HG2	1:A:58:ASP:N	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:ASN:OD1	1:A:228:ASP:HB2	2.21	0.41
1:A:255:ILE:HG23	1:A:256:ILE:N	2.36	0.41
1:A:411:MET:CE	1:B:442:LEU:CD2	2.98	0.41
1:A:484:TYR:CD1	1:A:484:TYR:C	2.99	0.41
1:A:503:TYR:CD2	1:A:504:GLU:N	2.89	0.41
1:B:133:GLN:HA	1:B:162:ILE:HB	2.03	0.41
1:B:137:LEU:CD2	1:B:168:ILE:HD11	2.37	0.41
1:B:404:MET:HE3	1:B:420:LEU:HD13	1.94	0.41
1:B:493:ASP:O	1:B:497:LYS:HG3	2.21	0.41
1:A:88:PHE:N	1:A:88:PHE:CD2	2.87	0.40
1:A:355:ILE:HD13	1:A:496:MET:HE3	2.03	0.40
1:B:32:HIS:CD2	1:B:62:PRO:HG2	2.56	0.40
1:A:211:LEU:CD2	1:A:392:ILE:CD1	2.99	0.40
1:B:240:PRO:HA	1:B:362:LEU:O	2.21	0.40
1:B:477:PRO:O	1:B:479:PRO:HD3	2.22	0.40
1:A:122:ASN:C	1:A:124:HIS:N	2.79	0.40
1:B:192:ARG:HG3	1:B:195:GLN:OE1	2.21	0.40
1:A:214:LEU:HD23	1:A:214:LEU:HA	1.84	0.40
1:A:323:LEU:C	1:A:325:ASP:H	2.29	0.40
1:B:307:GLN:N	1:B:480:ILE:O	2.50	0.40
1:B:405:MET:CE	1:B:416:GLU:O	2.67	0.40
1:B:63:GLU:O	1:B:104:ARG:NH2	2.55	0.40
1:B:176:ASP:N	1:B:179:SER:OG	2.49	0.40
1:B:275:LYS:HA	1:B:276:PRO:HD2	1.88	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	477/489 (98%)	417 (87%)	51 (11%)	9 (2%)	6 23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	477/489 (98%)	437 (92%)	35 (7%)	5 (1%)	12	39
All	All	954/978 (98%)	854 (90%)	86 (9%)	14 (2%)	8	28

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	B	158	CYS
1	B	504	GLU
1	A	73	SER
1	A	158	CYS
1	B	281	SER
1	A	123	SER
1	A	244	GLU
1	A	504	GLU
1	B	350	ASP
1	B	378	GLY
1	A	378	GLY
1	A	329	PRO
1	A	316	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	418/429 (97%)	351 (84%)	67 (16%)	2	8
1	B	419/429 (98%)	359 (86%)	60 (14%)	3	11
All	All	837/858 (98%)	710 (85%)	127 (15%)	3	9

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	VAL
1	A	28	GLN

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Mol	Chain	Res	Type
1	A	40	SER
1	A	43	LEU
1	A	47	LYS
1	A	48	ILE
1	A	53	TRP
1	A	57	ARG
1	A	80	ILE
1	A	82	LYS
1	A	83	GLN
1	A	94	GLU
1	A	104	ARG
1	A	106	SER
1	A	123	SER
1	A	132	SER
1	A	133	GLN
1	A	137	LEU
1	A	142	LEU
1	A	148	GLU
1	A	152	LYS
1	A	157	SER
1	A	160	SER
1	A	169	VAL
1	A	175	ARG
1	A	176	ASP
1	A	179	SER
1	A	180	SER
1	A	184	SER
1	A	192	ARG
1	A	199	ILE
1	A	212	MET
1	A	215	ARG
1	A	220	ILE
1	A	236	THR
1	A	238	LYS
1	A	256	ILE
1	A	260	MET
1	A	261	GLN
1	A	264	LEU
1	A	281	SER
1	A	283	ASP
1	A	285	ARG
1	A	288	LYS

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Mol	Chain	Res	Type
1	A	290	LYS
1	A	313	ASP
1	A	317	GLU
1	A	330	ARG
1	A	345	GLU
1	A	366	LYS
1	A	371	LEU
1	A	375	ASP
1	A	386	LYS
1	A	414	ASN
1	A	416	GLU
1	A	419	GLU
1	A	426	ASN
1	A	427	ARG
1	A	430	ASN
1	A	442	LEU
1	A	450	MET
1	A	456	ASP
1	A	459	ARG
1	A	463	ARG
1	A	478	LYS
1	A	497	LYS
1	A	502	GLN
1	B	29	SER
1	B	31	THR
1	B	35	ILE
1	B	37	MET
1	B	40	SER
1	B	73	SER
1	B	80	ILE
1	B	95	LYS
1	B	96	LEU
1	B	104	ARG
1	B	125	MET
1	B	132	SER
1	B	133	GLN
1	B	157	SER
1	B	159	MET
1	B	160	SER
1	B	170	GLU
1	B	177	LEU
1	B	179	SER

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Mol	Chain	Res	Type
1	B	182	ARG
1	B	184	SER
1	B	189	SER
1	B	190	LEU
1	B	212	MET
1	B	214	LEU
1	B	236	THR
1	B	255	ILE
1	B	260	MET
1	B	261	GLN
1	B	262	ASN
1	B	264	LEU
1	B	281	SER
1	B	283	ASP
1	B	285	ARG
1	B	299	GLN
1	B	301	ASN
1	B	305	LEU
1	B	330	ARG
1	B	332	SER
1	B	340	VAL
1	B	347	GLU
1	B	360	LYS
1	B	365	ARG
1	B	393	ARG
1	B	397	ASN
1	B	402	THR
1	B	404	MET
1	B	405	MET
1	B	414	ASN
1	B	419	GLU
1	B	426	ASN
1	B	427	ARG
1	B	432	LYS
1	B	441	ILE
1	B	470	HIS
1	B	473	GLU
1	B	476	LYS
1	B	491	GLU
1	B	498	ARG
1	B	504	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	155	HIS
1	A	261	GLN
1	A	311	ASN
1	A	374	HIS
1	B	32	HIS
1	B	226	ASN
1	B	229	ASN
1	B	261	GLN
1	B	299	GLN
1	B	382	HIS
1	B	451	HIS
1	B	471	GLN

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](#)

8 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	GOL	A	1506	-	5,5,5	0.53	0	5,5,5	0.68	0
3	BG6	B	1508	-	16,16,16	0.72	0	23,24,24	1.16	3 (13%)
2	GOL	A	1508	-	5,5,5	0.48	0	5,5,5	0.42	0
2	GOL	B	1506	-	5,5,5	0.36	0	5,5,5	0.46	0
2	GOL	B	1507	-	5,5,5	0.45	0	5,5,5	0.48	0
2	GOL	A	1509	-	5,5,5	0.36	0	5,5,5	0.41	0
3	BG6	A	1510	-	16,16,16	0.70	0	23,24,24	1.12	3 (13%)
2	GOL	A	1507	-	5,5,5	0.39	0	5,5,5	0.31	0

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	GOL	A	1506	-	-	1/4/4/4	-
3	BG6	B	1508	-	-	0/6/26/26	0/1/1/1
2	GOL	A	1508	-	-	0/4/4/4	-
2	GOL	B	1506	-	-	3/4/4/4	-
2	GOL	B	1507	-	-	3/4/4/4	-
2	GOL	A	1509	-	-	4/4/4/4	-
3	BG6	A	1510	-	-	1/6/26/26	0/1/1/1
2	GOL	A	1507	-	-	2/4/4/4	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1508	BG6	O2P-P-O6	-2.91	99.09	106.67
3	A	1510	BG6	O2P-P-O1P	2.72	118.02	107.80
3	B	1508	BG6	O2P-P-O1P	2.68	117.86	107.80
3	A	1510	BG6	O2P-P-O6	-2.65	99.75	106.67
3	A	1510	BG6	O1P-P-O6	-2.37	100.49	106.67
3	B	1508	BG6	O1P-P-O6	-2.20	100.94	106.67

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1509	GOL	O1-C1-C2-C3

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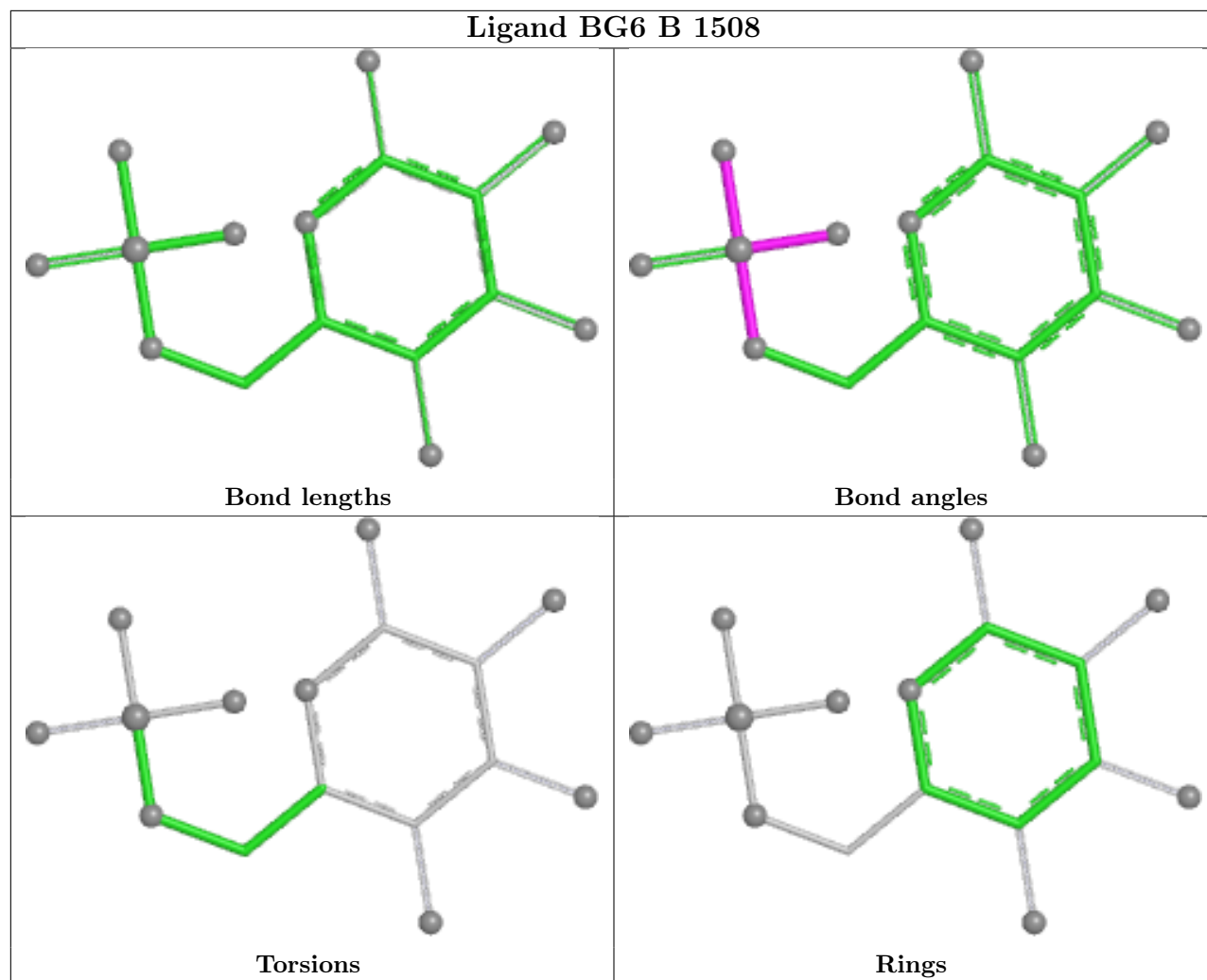
Mol	Chain	Res	Type	Atoms
2	A	1509	GOL	C1-C2-C3-O3
2	B	1506	GOL	O1-C1-C2-C3
2	B	1507	GOL	O1-C1-C2-C3
2	A	1509	GOL	O2-C2-C3-O3
2	A	1507	GOL	O1-C1-C2-C3
2	B	1506	GOL	C1-C2-C3-O3
2	B	1507	GOL	C1-C2-C3-O3
2	A	1509	GOL	O1-C1-C2-O2
2	B	1506	GOL	O1-C1-C2-O2
2	B	1507	GOL	O1-C1-C2-O2
2	A	1507	GOL	O1-C1-C2-O2
3	A	1510	BG6	O5-C5-C6-O6
2	A	1506	GOL	O2-C2-C3-O3

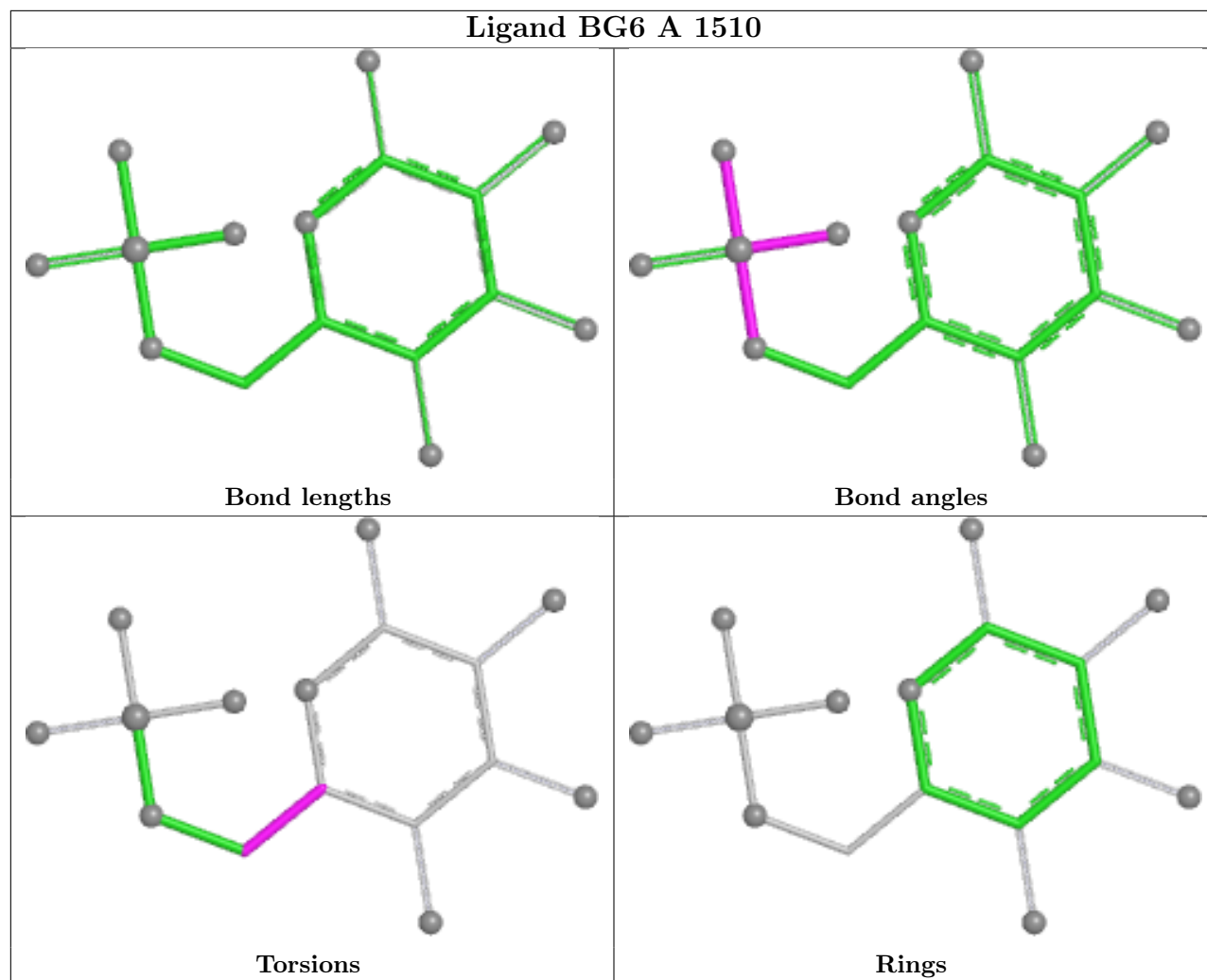
There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1506	GOL	2	0
3	B	1508	BG6	1	0
2	B	1506	GOL	5	0
2	A	1507	GOL	7	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

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	479/489 (97%)	0.44	23 (4%) 35 28	25, 47, 72, 94	1 (0%)
1	B	479/489 (97%)	0.50	20 (4%) 40 32	30, 51, 76, 94	0
All	All	958/978 (97%)	0.47	43 (4%) 38 30	25, 49, 74, 94	1 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	504	GLU	10.5
1	B	27	VAL	4.4
1	B	505	GLY	4.0
1	A	89	LYS	3.9
1	A	118	TYR	3.8
1	B	503	TYR	3.7
1	B	243	THR	3.6
1	A	157	SER	3.3
1	A	27	VAL	3.1
1	A	176	ASP	3.1
1	B	176	ASP	3.0
1	B	385	CYS	3.0
1	A	243	THR	2.9
1	B	245	GLY	2.8
1	B	170	GLU	2.8
1	A	143	PRO	2.8
1	A	503	TYR	2.7
1	A	249	TYR	2.7
1	A	505	GLY	2.7
1	B	317	GLU	2.6
1	A	160	SER	2.6
1	A	108	VAL	2.6
1	B	253	PHE	2.5
1	B	330	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	122	ASN	2.4
1	A	146	VAL	2.3
1	A	90	ALA	2.3
1	B	314	GLY	2.2
1	A	245	GLY	2.2
1	B	312	PRO	2.1
1	B	228	ASP	2.1
1	A	140	LEU	2.1
1	A	156	GLU	2.1
1	A	252	GLU	2.1
1	A	247	GLY	2.1
1	A	158	CYS	2.1
1	B	180	SER	2.1
1	B	474	LEU	2.1
1	B	402	THR	2.1
1	A	42	ASP	2.0
1	B	171	LYS	2.0
1	B	244	GLU	2.0
1	B	320	LYS	2.0

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 oligosaccharides 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.

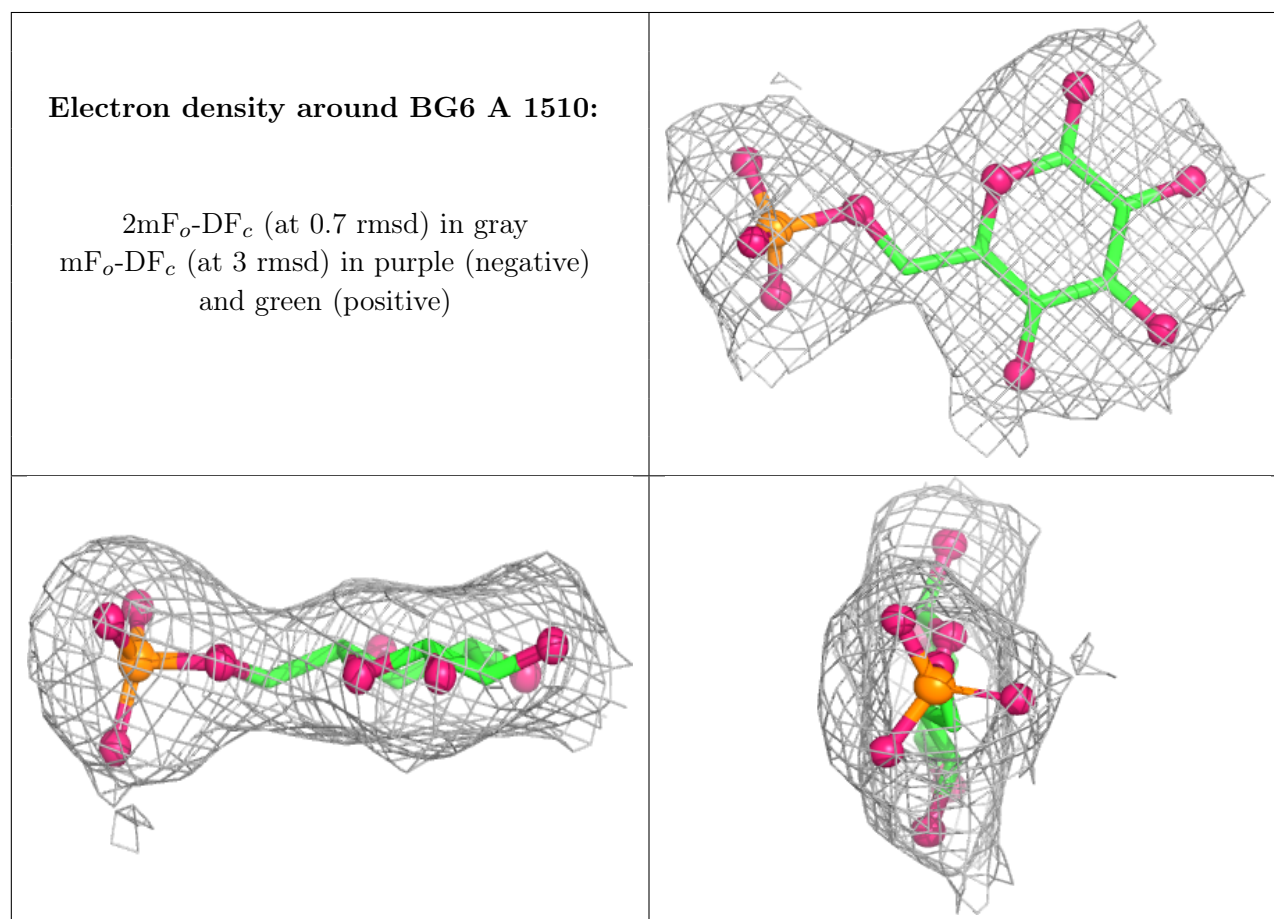
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GOL	B	1507	6/6	0.62	0.24	53,53,53,53	0
2	GOL	B	1506	6/6	0.77	0.20	58,58,59,59	0
2	GOL	A	1509	6/6	0.81	0.20	59,59,59,60	0
2	GOL	A	1506	6/6	0.85	0.23	50,51,52,53	0

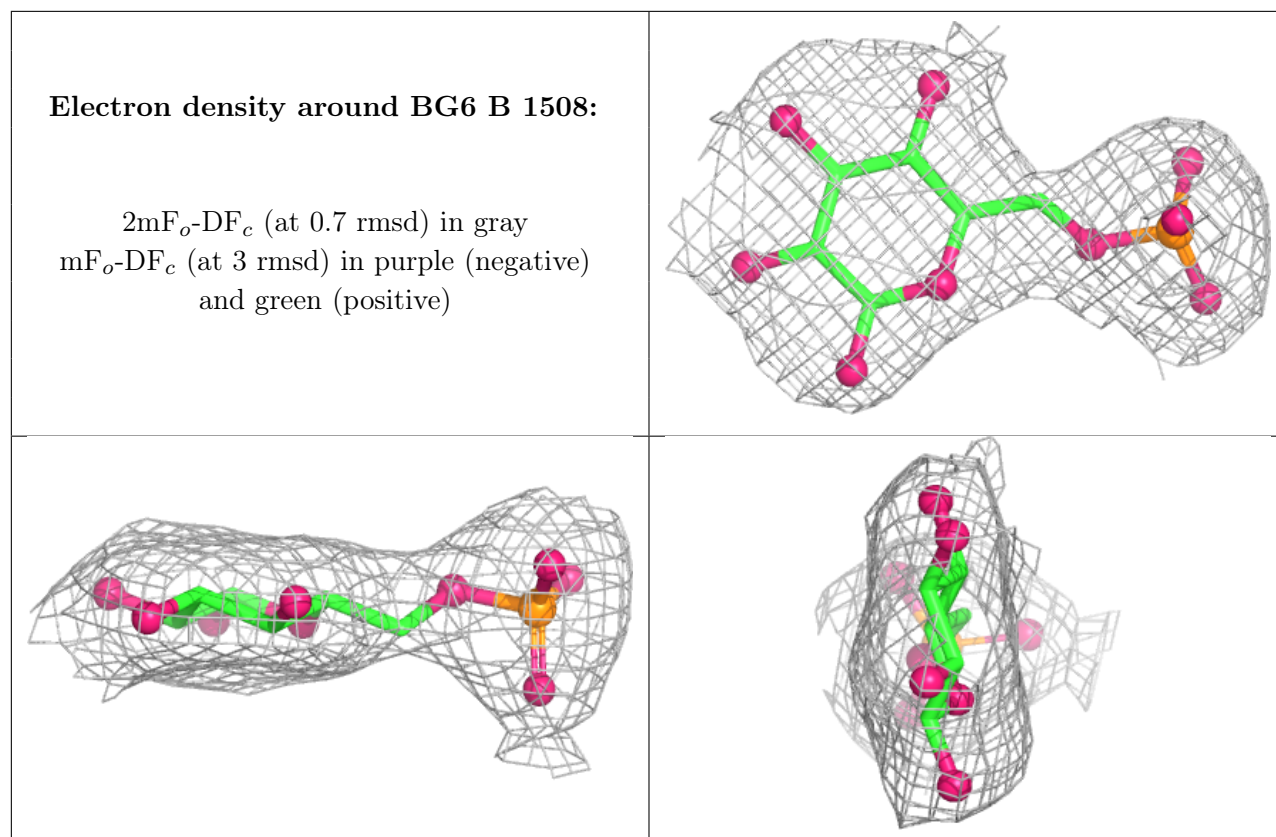
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	A	1507	6/6	0.87	0.14	52,52,52,53	0
2	GOL	A	1508	6/6	0.89	0.15	50,51,52,52	0
3	BG6	A	1510	16/16	0.94	0.09	45,46,48,48	0
3	BG6	B	1508	16/16	0.96	0.07	50,50,51,51	0

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.





6.5 Other polymers [i](#)

There are no such residues in this entry.