



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:55 PM UTC

PDB ID : 1DQR / pdb_00001dqr
Title : CRYSTAL STRUCTURE OF RABBIT PHOSPHOGLUCOSE ISOMERASE,
A GLYCOLYTIC ENZYME THAT MOONLIGHTS AS NEUROLEUKIN,
AUTOCRINE MOTILITY FACTOR, AND DIFFERENTIATION MEDIA-
TOR
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Deposited on : 2000-01-05
Resolution : 2.50 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](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (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

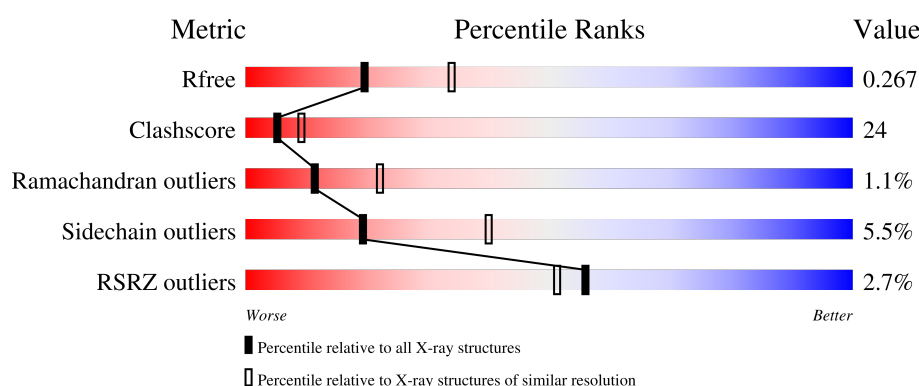
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	557	3% 60% 35% .
1	B	557	2% 58% 35% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9194 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.

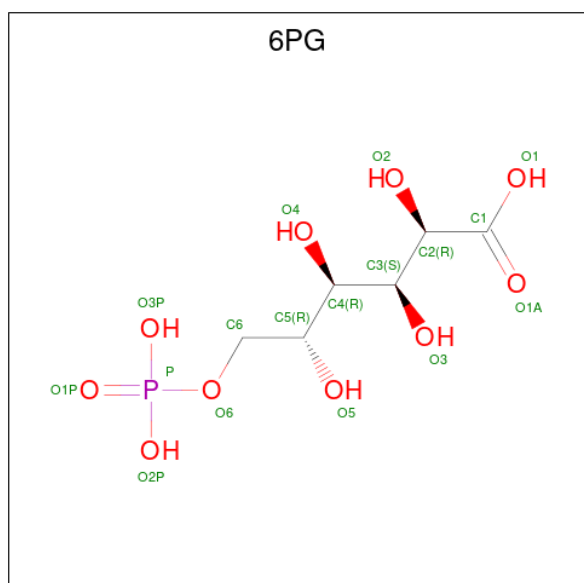
- Molecule 1 is a protein called PHOSPHOGLUCOSE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4395	2805	765	806	19			
1	B	555	Total	C	N	O	S	0	0	0
			4395	2805	765	806	19			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	109	PRO	LEU	see remark 999	UNP Q9N1E2
A	222	LYS	GLU	see remark 999	UNP Q9N1E2
A	283	ILE	VAL	see remark 999	UNP Q9N1E2
B	109	PRO	LEU	see remark 999	UNP Q9N1E2
B	222	LYS	GLU	see remark 999	UNP Q9N1E2
B	283	ILE	VAL	see remark 999	UNP Q9N1E2

- Molecule 2 is 6-PHOSPHOGLUCONIC ACID (CCD ID: 6PG) (formula: $C_6H_{13}O_{10}P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			17	6	10	1		
2	B	1	Total	C	O	P	0	0
			17	6	10	1		

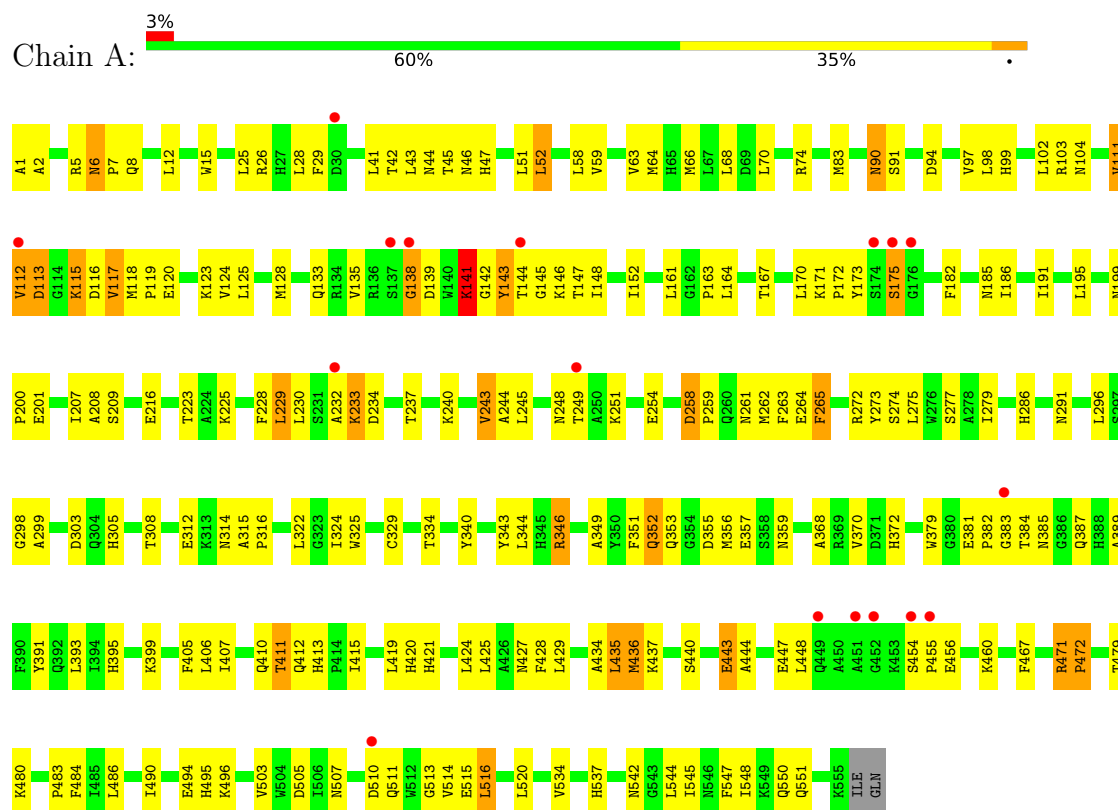
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	176	Total	O	0	0
			176	176		
3	B	194	Total	O	0	0
			194	194		

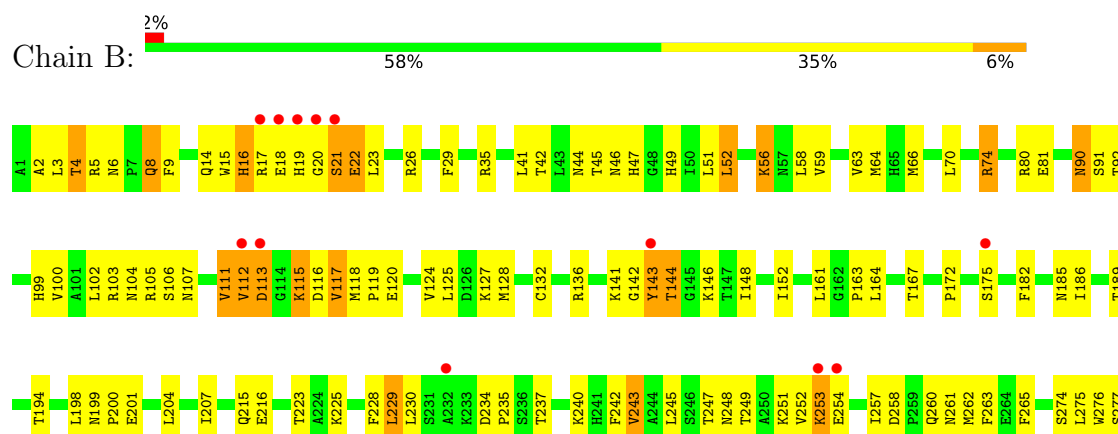
3 Residue-property plots

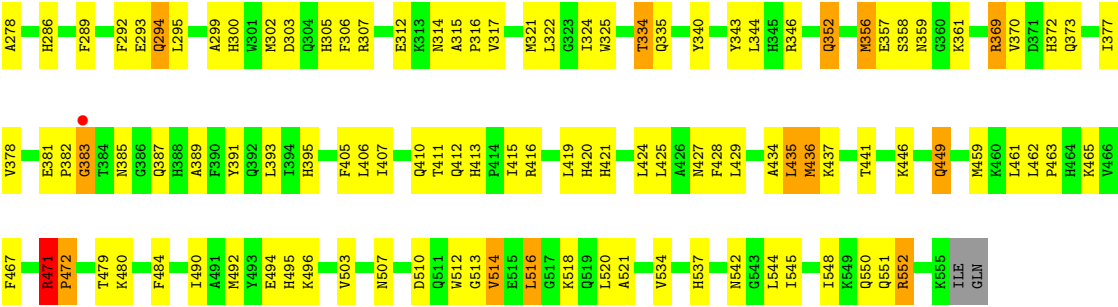
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: PHOSPHOGLUCOSE ISOMERASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	82.69Å 115.27Å 271.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.00 – 2.50 35.00 – 2.51	Depositor EDS
% Data completeness (in resolution range)	90.9 (35.00-2.50) 92.0 (35.00-2.51)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.48 (at 2.50Å)	Xtriage
Refinement program	CNS, XTALVIEW	Depositor
R, R_{free}	0.212 , 0.265 0.213 , 0.267	Depositor DCC
R_{free} test set	4152 reflections (10.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9194	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: 6PG

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	0.47	0/4504	1.02	23/6101 (0.4%)
1	B	0.48	1/4504 (0.0%)	0.99	20/6101 (0.3%)
All	All	0.48	1/9008 (0.0%)	1.01	43/12202 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	514	VAL	CB-CG1	-5.18	1.35	1.52

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	LYS	N-CA-C	12.06	129.22	107.99
1	A	352	GLN	N-CA-C	-9.69	100.71	111.07
1	B	352	GLN	N-CA-C	-8.96	101.51	111.28
1	A	175	SER	N-CA-C	8.62	129.16	110.80
1	B	175	SER	N-CA-C	8.18	128.23	110.80
1	B	471	ARG	CA-C-N	8.13	128.61	119.83
1	B	471	ARG	C-N-CA	8.13	128.61	119.83
1	B	112	VAL	CB-CA-C	-7.91	98.31	111.29
1	A	112	VAL	CB-CA-C	-7.76	98.56	111.29
1	A	471	ARG	CA-C-N	7.75	128.20	119.83
1	A	471	ARG	C-N-CA	7.75	128.20	119.83
1	A	6	ASN	CA-C-N	7.38	127.24	119.05
1	A	6	ASN	C-N-CA	7.38	127.24	119.05
1	A	359	ASN	N-CA-C	7.07	120.85	112.93
1	A	98	LEU	N-CA-C	7.04	119.18	109.54
1	A	111	VAL	N-CA-C	7.01	118.38	108.36
1	A	258	ASP	CA-C-N	6.67	128.17	119.84
1	A	258	ASP	C-N-CA	6.67	128.17	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	334	THR	N-CA-C	6.62	120.30	109.85
1	B	334	THR	N-CA-C	6.47	120.08	109.85
1	A	91	SER	N-CA-C	6.27	118.67	111.02
1	B	449	GLN	N-CA-C	-6.22	103.81	111.40
1	B	91	SER	N-CA-C	6.21	118.56	111.11
1	B	111	VAL	N-CA-C	6.21	117.24	108.36
1	A	496	LYS	N-CA-C	-6.13	104.51	111.07
1	B	247	THR	N-CA-C	-5.94	104.51	113.89
1	B	496	LYS	N-CA-C	-5.85	104.90	111.28
1	A	308	THR	N-CA-C	5.68	120.32	113.17
1	B	21	SER	N-CA-C	-5.68	106.74	113.38
1	B	307	ARG	N-CA-C	5.66	117.53	111.36
1	B	514	VAL	CB-CA-C	-5.57	105.14	111.55
1	B	294	GLN	N-CA-C	-5.49	105.45	111.82
1	A	443	GLU	N-CA-C	-5.48	105.31	111.28
1	B	4	THR	N-CA-C	-5.42	106.62	113.18
1	A	274	SER	N-CA-C	5.24	119.72	113.23
1	B	215	GLN	N-CA-C	5.22	116.97	111.28
1	B	274	SER	N-CA-C	5.20	119.78	113.38
1	B	356	MET	N-CA-C	5.16	117.31	111.11
1	A	472	PRO	N-CA-C	5.13	118.92	111.03
1	A	244	ALA	N-CA-C	5.11	117.14	109.23
1	A	515	GLU	N-CA-C	5.05	117.67	111.82
1	B	472	PRO	N-CA-C	5.04	118.79	111.03
1	A	264	GLU	N-CA-C	5.03	118.22	110.42

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	4395	0	4352	227	0
1	B	4395	0	4352	229	0
2	A	17	0	10	1	0
2	B	17	0	10	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	176	0	0	12	0
3	B	194	0	0	16	0
All	All	9194	0	8724	428	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 24.

All (428) 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:143:TYR:CE2	3:B:787:HOH:O	1.87	1.24
1:A:113:ASP:CG	3:A:724:HOH:O	1.89	1.12
1:A:1:ALA:CB	3:A:774:HOH:O	1.95	1.10
1:B:112:VAL:HG23	1:B:112:VAL:O	1.48	1.10
1:A:112:VAL:HG23	1:A:112:VAL:O	1.38	1.08
1:B:112:VAL:O	1:B:113:ASP:HB2	1.61	1.00
1:A:142:GLY:O	1:A:145:GLY:N	1.95	0.98
1:A:124:VAL:HG12	1:A:128:MET:HE2	1.50	0.93
1:A:141:LYS:HB2	1:A:240:LYS:HE2	1.50	0.93
1:B:550:GLN:HE21	1:B:551:GLN:HE21	1.09	0.92
1:B:124:VAL:HG12	1:B:128:MET:HE2	1.51	0.92
1:B:143:TYR:CD2	3:B:787:HOH:O	2.13	0.91
1:A:112:VAL:O	1:A:112:VAL:CG2	2.13	0.91
1:A:1:ALA:HB1	3:A:774:HOH:O	1.61	0.91
1:B:6:ASN:HD21	1:B:8:GLN:HG2	1.37	0.89
1:B:240:LYS:HE3	3:B:787:HOH:O	1.76	0.86
1:A:125:LEU:HA	1:A:128:MET:HE3	1.56	0.85
1:B:90:ASN:ND2	1:B:507:ASN:HD21	1.78	0.81
1:A:59:VAL:HG23	1:A:63:VAL:HB	1.62	0.81
1:A:185:ASN:HD21	1:B:383:GLY:HA3	1.46	0.81
1:B:59:VAL:HG23	1:B:63:VAL:HB	1.61	0.81
1:B:125:LEU:HA	1:B:128:MET:HE3	1.61	0.79
1:A:5:ARG:HH11	1:A:5:ARG:HB2	1.48	0.79
1:A:272:ARG:HH12	1:A:511:GLN:HE21	1.32	0.77
1:B:550:GLN:HE21	1:B:551:GLN:NE2	1.82	0.77
1:A:112:VAL:O	1:A:113:ASP:HB2	1.85	0.77
1:B:143:TYR:HE2	3:B:787:HOH:O	1.33	0.77
1:A:1:ALA:N	1:A:368:ALA:HB1	2.00	0.76
1:A:5:ARG:HB2	1:A:5:ARG:NH1	2.01	0.76
1:A:90:ASN:ND2	1:A:507:ASN:HD21	1.84	0.75
1:B:103:ARG:HD2	1:B:303:ASP:OD2	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:VAL:HG21	1:B:64:MET:SD	2.25	0.75
1:B:258:ASP:OD2	1:B:260:GLN:HG2	1.87	0.75
1:B:471:ARG:NH2	3:B:607:HOH:O	2.19	0.75
1:B:14:GLN:O	1:B:18:GLU:HB2	1.87	0.74
1:A:383:GLY:HA3	1:B:185:ASN:HD21	1.52	0.74
1:A:139:ASP:O	1:A:139:ASP:OD2	2.04	0.73
1:A:142:GLY:O	1:A:144:THR:N	2.21	0.73
1:A:209:SER:HA	2:A:600:6PG:O2P	1.87	0.73
1:A:44:ASN:ND2	1:A:46:ASN:H	1.87	0.73
1:A:1:ALA:H1	1:A:368:ALA:HB1	1.55	0.72
1:B:102:LEU:HD13	1:B:277:SER:HB3	1.72	0.72
1:A:185:ASN:ND2	1:B:383:GLY:HA3	2.03	0.72
1:B:6:ASN:HD22	1:B:9:PHE:H	1.38	0.71
1:B:550:GLN:NE2	1:B:551:GLN:HE21	1.87	0.71
1:B:248:ASN:ND2	1:B:251:LYS:HG3	2.05	0.71
1:B:315:ALA:HB3	1:B:316:PRO:HD3	1.71	0.71
1:A:353:GLN:O	1:A:357:GLU:HB2	1.91	0.71
1:B:324:ILE:HD13	1:B:503:VAL:HG11	1.73	0.70
1:A:315:ALA:HB3	1:A:316:PRO:HD3	1.73	0.70
1:A:172:PRO:HG2	1:A:412:GLN:HG3	1.72	0.70
1:A:550:GLN:HE21	1:A:551:GLN:HE21	1.39	0.70
1:B:240:LYS:HG3	3:B:787:HOH:O	1.92	0.70
1:B:253:LYS:HA	1:B:253:LYS:HE2	1.71	0.70
1:A:324:ILE:HD13	1:A:503:VAL:HG11	1.73	0.70
1:A:436:MET:HE2	1:A:437:LYS:HA	1.73	0.69
1:B:112:VAL:O	1:B:112:VAL:CG2	2.21	0.69
1:A:94:ASP:OD1	3:A:624:HOH:O	2.11	0.69
1:A:59:VAL:HG21	1:A:64:MET:SD	2.32	0.68
1:B:74:ARG:HD3	1:B:503:VAL:O	1.92	0.68
1:B:172:PRO:HG2	1:B:412:GLN:HG3	1.76	0.68
1:A:102:LEU:HD13	1:A:277:SER:HB3	1.75	0.68
1:A:383:GLY:HA3	1:B:185:ASN:ND2	2.09	0.68
1:A:349:ALA:HA	1:A:352:GLN:HG3	1.76	0.67
1:A:141:LYS:HG3	1:A:240:LYS:CE	2.24	0.67
1:A:407:ILE:HD13	1:A:425:LEU:HD23	1.76	0.67
1:A:340:TYR:HB3	1:A:424:LEU:HD11	1.76	0.67
1:B:436:MET:HE2	1:B:436:MET:C	2.20	0.67
3:A:765:HOH:O	1:B:49:HIS:HD2	1.77	0.66
1:B:436:MET:HE2	1:B:437:LYS:HA	1.75	0.66
1:A:8:GLN:H	1:A:8:GLN:NE2	1.94	0.66
1:B:254:GLU:HG3	3:B:671:HOH:O	1.93	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:537:HIS:H	1:A:542:ASN:HD21	1.43	0.65
1:B:59:VAL:HG23	1:B:63:VAL:CB	2.27	0.65
1:B:537:HIS:H	1:B:542:ASN:HD21	1.42	0.65
1:A:143:TYR:CZ	1:A:232:ALA:HB2	2.32	0.65
1:A:146:LYS:HD3	1:A:147:THR:H	1.63	0.64
1:B:17:ARG:NH1	1:B:17:ARG:HB2	2.12	0.64
1:B:361:LYS:HZ3	1:B:510:ASP:HB2	1.61	0.64
1:B:441:THR:HG23	1:B:462:LEU:HD11	1.78	0.64
1:B:148:ILE:HD13	1:B:204:LEU:HB2	1.78	0.64
1:A:243:VAL:HG22	1:A:261:ASN:O	1.99	0.63
1:B:387:GLN:HE22	1:B:427:ASN:HB3	1.63	0.63
1:A:59:VAL:HG23	1:A:63:VAL:CB	2.28	0.63
1:B:343:TYR:OH	1:B:413:HIS:HE1	1.82	0.63
1:A:141:LYS:HE3	1:A:141:LYS:N	2.13	0.62
1:B:357:GLU:OE2	2:B:601:6PG:O1	2.17	0.62
1:A:436:MET:HE2	1:A:436:MET:C	2.25	0.62
1:B:141:LYS:O	1:B:240:LYS:HD2	1.98	0.62
1:A:387:GLN:HE22	1:A:427:ASN:HB3	1.65	0.61
1:B:340:TYR:HB3	1:B:424:LEU:HD11	1.82	0.61
1:A:436:MET:HE2	1:A:437:LYS:CA	2.30	0.61
1:B:411:THR:HG23	1:B:421:HIS:CE1	2.34	0.61
1:A:314:ASN:OD1	1:A:316:PRO:HD2	2.01	0.61
1:A:385:ASN:H	1:A:385:ASN:HD22	1.48	0.61
1:B:103:ARG:HD3	1:B:299:ALA:CB	2.31	0.61
1:B:436:MET:HE2	1:B:437:LYS:CA	2.31	0.61
1:A:141:LYS:CB	1:A:240:LYS:HE2	2.27	0.61
1:A:411:THR:HG23	1:A:421:HIS:CE1	2.36	0.61
1:B:436:MET:HE2	1:B:437:LYS:N	2.15	0.60
1:B:59:VAL:CG2	1:B:63:VAL:HB	2.32	0.60
1:A:144:THR:CG2	1:A:201:GLU:HB2	2.31	0.60
1:B:44:ASN:ND2	1:B:46:ASN:H	1.99	0.60
1:B:59:VAL:HG23	1:B:63:VAL:CG1	2.31	0.60
1:A:233:LYS:HB2	1:A:233:LYS:HZ3	1.67	0.60
1:B:385:ASN:HD22	1:B:385:ASN:H	1.47	0.60
1:B:20:GLY:C	1:B:22:GLU:H	2.10	0.60
1:A:436:MET:HE2	1:A:437:LYS:N	2.17	0.60
1:A:343:TYR:OH	1:A:413:HIS:HE1	1.85	0.59
1:B:395:HIS:HD2	3:B:660:HOH:O	1.84	0.59
1:A:58:LEU:CD2	1:A:472:PRO:HB3	2.32	0.59
1:A:103:ARG:NH2	1:A:494:GLU:OE1	2.35	0.59
1:A:59:VAL:HG23	1:A:63:VAL:CG1	2.33	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:VAL:CG2	1:A:63:VAL:HB	2.31	0.59
1:B:42:THR:HG23	3:B:632:HOH:O	2.02	0.59
1:A:233:LYS:HB2	1:A:233:LYS:NZ	2.18	0.59
1:B:407:ILE:HD13	1:B:425:LEU:HD23	1.83	0.59
1:A:141:LYS:N	1:A:141:LYS:CE	2.65	0.58
1:B:352:GLN:HG3	1:B:378:VAL:O	2.03	0.58
1:B:544:LEU:O	1:B:548:ILE:HG13	2.03	0.58
1:A:381:GLU:HG3	1:A:382:PRO:HD2	1.84	0.58
1:B:275:LEU:HD12	1:B:275:LEU:C	2.29	0.58
1:A:142:GLY:C	1:A:144:THR:N	2.60	0.58
1:B:143:TYR:CD1	1:B:237:THR:HG22	2.38	0.58
1:A:249:THR:HA	1:A:262:MET:SD	2.44	0.57
1:B:58:LEU:CD2	1:B:472:PRO:HB3	2.35	0.57
1:B:146:LYS:NZ	1:B:201:GLU:OE1	2.36	0.57
1:A:312:GLU:H	1:A:312:GLU:CD	2.12	0.57
1:B:302:MET:HE3	1:B:492:MET:HB2	1.87	0.57
1:B:361:LYS:NZ	1:B:510:ASP:HB2	2.19	0.56
1:A:410:GLN:HG3	1:A:480:LYS:HD2	1.87	0.56
1:B:234:ASP:HB3	1:B:237:THR:HG23	1.85	0.56
1:B:257:ILE:HG22	1:B:258:ASP:N	2.21	0.56
1:A:234:ASP:HB3	1:A:237:THR:HG23	1.86	0.56
1:A:141:LYS:O	1:A:240:LYS:HB3	2.04	0.56
1:A:117:VAL:O	1:A:120:GLU:HG2	2.06	0.56
1:B:117:VAL:O	1:B:120:GLU:HG2	2.06	0.56
1:A:90:ASN:C	1:A:90:ASN:HD22	2.12	0.56
1:B:253:LYS:HE2	1:B:253:LYS:CA	2.34	0.56
1:B:164:LEU:C	1:B:164:LEU:HD23	2.31	0.56
1:B:6:ASN:ND2	1:B:8:GLN:HG2	2.13	0.56
1:B:56:LYS:HB2	3:B:621:HOH:O	2.06	0.55
1:A:443:GLU:O	1:A:447:GLU:HG3	2.07	0.55
1:A:275:LEU:C	1:A:275:LEU:HD12	2.31	0.55
1:B:490:ILE:O	1:B:494:GLU:HG3	2.06	0.55
1:A:103:ARG:HD3	1:A:299:ALA:CB	2.37	0.55
1:A:420:HIS:HD2	1:B:186:ILE:O	1.90	0.55
1:B:100:VAL:HG22	3:B:617:HOH:O	2.06	0.55
1:A:103:ARG:HD2	1:A:303:ASP:OD2	2.07	0.55
1:A:186:ILE:HD13	1:A:216:GLU:HA	1.88	0.55
1:A:353:GLN:HA	1:B:389:ALA:O	2.06	0.55
1:A:223:THR:OG1	1:B:420:HIS:HE1	1.90	0.54
1:B:393:LEU:HD23	1:B:393:LEU:O	2.07	0.54
1:B:462:LEU:C	1:B:462:LEU:HD13	2.32	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:VAL:HG12	1:A:372:HIS:H	1.71	0.54
1:B:167:THR:HG23	1:B:286:HIS:ND1	2.23	0.54
1:A:141:LYS:CD	1:A:141:LYS:H	2.19	0.54
1:A:490:ILE:O	1:A:494:GLU:HG3	2.08	0.54
1:A:135:VAL:HG13	1:A:148:ILE:CG1	2.38	0.54
1:A:185:ASN:HD21	1:B:383:GLY:CA	2.19	0.54
1:A:26:ARG:HA	1:A:436:MET:HE3	1.90	0.53
1:B:446:LYS:HE3	1:B:449:GLN:HE22	1.73	0.53
1:A:545:ILE:HG12	1:B:429:LEU:HD13	1.89	0.53
1:B:90:ASN:C	1:B:90:ASN:HD22	2.15	0.53
1:B:312:GLU:CD	1:B:312:GLU:H	2.16	0.53
1:B:199:ASN:OD1	1:B:201:GLU:HG2	2.09	0.53
1:A:186:ILE:CG2	1:B:424:LEU:HD23	2.39	0.53
1:A:393:LEU:O	1:A:393:LEU:HD23	2.09	0.53
1:B:26:ARG:CA	1:B:436:MET:HE3	2.39	0.53
1:A:15:TRP:CG	1:A:66:MET:HE1	2.44	0.53
1:B:152:ILE:N	1:B:152:ILE:HD12	2.24	0.53
1:A:8:GLN:H	1:A:8:GLN:HE21	1.56	0.52
1:A:113:ASP:HB2	3:A:724:HOH:O	2.08	0.52
1:A:164:LEU:C	1:A:164:LEU:HD23	2.35	0.52
1:A:135:VAL:HG13	1:A:148:ILE:HG12	1.90	0.52
1:A:207:ILE:HD12	1:A:207:ILE:N	2.25	0.52
1:B:45:THR:O	1:B:46:ASN:HB2	2.09	0.52
1:B:381:GLU:HG3	1:B:382:PRO:HD2	1.90	0.52
1:A:167:THR:HG23	1:A:286:HIS:ND1	2.24	0.52
1:A:164:LEU:HD22	1:A:346:ARG:HG2	1.91	0.52
1:B:370:VAL:HG12	1:B:372:HIS:H	1.75	0.52
1:B:26:ARG:HA	1:B:436:MET:HE3	1.90	0.52
1:B:163:PRO:HB2	1:B:182:PHE:HE2	1.75	0.52
1:B:20:GLY:C	1:B:22:GLU:N	2.63	0.52
1:B:143:TYR:HD1	1:B:237:THR:HG22	1.75	0.52
1:A:142:GLY:O	1:A:143:TYR:C	2.53	0.51
1:A:199:ASN:OD1	1:A:201:GLU:HG2	2.10	0.51
1:A:415:ILE:HD11	1:A:420:HIS:CG	2.45	0.51
1:B:99:HIS:HE1	3:B:604:HOH:O	1.91	0.51
1:A:410:GLN:HG2	1:A:479:THR:O	2.10	0.51
1:A:44:ASN:HD22	1:A:45:THR:N	2.09	0.51
1:B:186:ILE:HD13	1:B:216:GLU:HA	1.91	0.51
1:A:303:ASP:OD1	1:A:495:HIS:CE1	2.63	0.51
1:B:141:LYS:O	1:B:141:LYS:HD2	2.10	0.51
1:B:207:ILE:N	1:B:207:ILE:HD12	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ALA:H2	1:A:368:ALA:HB1	1.76	0.51
1:B:80:ARG:HG3	1:B:306:PHE:CE2	2.46	0.51
1:B:243:VAL:HG22	1:B:261:ASN:O	2.11	0.51
1:A:223:THR:OG1	1:B:416:ARG:NH1	2.43	0.51
1:A:116:ASP:OD2	1:A:119:PRO:HD3	2.11	0.51
1:B:21:SER:O	1:B:22:GLU:C	2.53	0.51
1:B:245:LEU:HA	1:B:263:PHE:O	2.11	0.51
1:B:305:HIS:HE1	1:B:315:ALA:H	1.58	0.50
1:A:152:ILE:HD12	1:A:152:ILE:N	2.26	0.50
1:A:395:HIS:NE2	1:A:435:LEU:HD13	2.27	0.50
1:B:314:ASN:OD1	1:B:316:PRO:HD2	2.11	0.50
1:B:41:LEU:HD12	1:B:41:LEU:O	2.11	0.50
1:B:99:HIS:CE1	3:B:604:HOH:O	2.64	0.50
1:B:471:ARG:HB3	3:B:744:HOH:O	2.12	0.50
1:B:248:ASN:HD21	1:B:251:LYS:HG3	1.77	0.50
1:A:186:ILE:HG23	1:B:424:LEU:HD23	1.94	0.50
1:B:141:LYS:HD2	1:B:141:LYS:C	2.37	0.50
1:A:113:ASP:CB	3:A:724:HOH:O	2.42	0.50
1:A:249:THR:HG22	1:A:262:MET:SD	2.51	0.50
1:B:393:LEU:HD23	1:B:393:LEU:C	2.37	0.50
1:A:272:ARG:HH12	1:A:511:GLN:NE2	2.05	0.49
1:B:415:ILE:HD12	1:B:415:ILE:C	2.37	0.49
1:A:2:ALA:HA	1:A:5:ARG:HD3	1.94	0.49
1:A:141:LYS:HG3	1:A:240:LYS:HD2	1.94	0.49
1:A:142:GLY:C	1:A:144:THR:H	2.20	0.49
1:A:429:LEU:HD13	1:B:545:ILE:HG12	1.94	0.49
1:A:440:SER:OG	1:A:443:GLU:HG3	2.12	0.49
1:A:163:PRO:HB2	1:A:182:PHE:HE2	1.77	0.49
1:B:303:ASP:OD1	1:B:495:HIS:HE1	1.96	0.49
1:B:410:GLN:HE21	1:B:480:LYS:HD2	1.76	0.49
1:B:14:GLN:O	1:B:18:GLU:N	2.43	0.49
1:A:104:ASN:HD22	1:A:118:MET:HE2	1.78	0.49
1:A:393:LEU:HD23	1:A:393:LEU:C	2.38	0.49
1:B:6:ASN:HD22	1:B:9:PHE:N	2.08	0.49
1:B:125:LEU:HD23	1:B:128:MET:HE3	1.95	0.49
1:A:45:THR:O	1:A:46:ASN:HB2	2.13	0.48
1:B:248:ASN:O	1:B:252:VAL:HG23	2.12	0.48
1:B:292:PHE:O	1:B:295:LEU:HB3	2.12	0.48
1:B:461:LEU:HG	1:B:465:LYS:HD2	1.94	0.48
1:A:208:ALA:HB2	1:A:245:LEU:HD12	1.94	0.48
1:B:15:TRP:CG	1:B:66:MET:HE1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:LEU:HD12	1:A:41:LEU:O	2.13	0.48
1:B:411:THR:HG23	1:B:421:HIS:ND1	2.29	0.48
1:A:138:GLY:HA2	1:A:147:THR:HG22	1.94	0.48
1:A:513:GLY:HA2	1:B:467:PHE:CZ	2.49	0.48
1:B:534:VAL:HG21	1:B:545:ILE:HG21	1.94	0.48
1:A:143:TYR:CE2	1:A:232:ALA:HB2	2.48	0.48
1:A:413:HIS:ND1	1:B:189:THR:HG22	2.29	0.48
1:A:26:ARG:CA	1:A:436:MET:HE3	2.43	0.48
1:A:424:LEU:HD23	1:B:186:ILE:CG2	2.44	0.48
1:B:17:ARG:C	1:B:19:HIS:N	2.72	0.48
1:A:415:ILE:C	1:A:415:ILE:HD12	2.39	0.48
1:B:289:PHE:O	1:B:293:GLU:HG3	2.14	0.47
1:A:186:ILE:HB	1:A:216:GLU:HG3	1.95	0.47
1:A:467:PHE:CZ	1:B:513:GLY:HA2	2.50	0.47
1:B:518:LYS:O	1:B:521:ALA:HB3	2.14	0.47
1:B:242:PHE:HB3	1:B:257:ILE:CD1	2.45	0.47
1:B:103:ARG:HD3	1:B:299:ALA:HB3	1.94	0.47
1:A:434:ALA:HB2	1:B:520:LEU:HD12	1.97	0.47
1:A:510:ASP:HB3	3:A:767:HOH:O	2.15	0.47
1:B:132:CYS:O	1:B:136:ARG:HG3	2.14	0.47
1:B:510:ASP:HB3	3:B:784:HOH:O	2.14	0.47
1:A:12:LEU:HD23	1:A:66:MET:CB	2.45	0.47
1:A:104:ASN:HD22	1:A:118:MET:CE	2.28	0.47
1:B:41:LEU:HD12	1:B:41:LEU:C	2.40	0.47
1:B:103:ARG:HD3	1:B:299:ALA:HB1	1.96	0.47
1:B:358:SER:C	1:B:359:ASN:HD22	2.22	0.47
1:B:410:GLN:HG2	1:B:479:THR:O	2.13	0.47
1:B:462:LEU:HD13	1:B:463:PRO:N	2.28	0.47
1:A:272:ARG:NH1	1:A:511:GLN:HE21	2.07	0.47
1:A:411:THR:HG23	1:A:421:HIS:ND1	2.30	0.47
1:B:303:ASP:OD1	1:B:495:HIS:CE1	2.68	0.47
1:A:41:LEU:HD12	1:A:41:LEU:C	2.39	0.47
1:A:104:ASN:ND2	1:A:118:MET:HE2	2.30	0.47
1:B:294:GLN:O	1:B:484:PHE:HA	2.14	0.47
1:B:103:ARG:O	1:B:300:HIS:HB2	2.14	0.47
1:B:361:LYS:NZ	1:B:510:ASP:CB	2.78	0.47
1:B:395:HIS:NE2	1:B:435:LEU:HD13	2.30	0.47
1:A:125:LEU:HD23	1:A:128:MET:HE3	1.97	0.46
1:B:52:LEU:C	1:B:52:LEU:CD2	2.88	0.46
1:B:142:GLY:O	1:B:144:THR:N	2.39	0.46
1:A:144:THR:HB	1:A:201:GLU:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:THR:HG22	1:B:51:LEU:HD13	1.96	0.46
1:B:17:ARG:O	1:B:17:ARG:HG2	2.14	0.46
1:A:355:ASP:HB2	3:A:746:HOH:O	2.14	0.46
1:A:544:LEU:O	1:A:548:ILE:HG13	2.15	0.46
1:B:128:MET:SD	1:B:278:ALA:HA	2.55	0.46
1:B:243:VAL:HG23	1:B:261:ASN:HB3	1.97	0.46
1:A:90:ASN:ND2	1:A:90:ASN:C	2.74	0.46
1:A:111:VAL:HG13	1:A:115:LYS:N	2.31	0.46
1:A:385:ASN:H	1:A:385:ASN:ND2	2.14	0.46
1:A:387:GLN:HA	1:A:391:TYR:CG	2.51	0.46
1:B:14:GLN:O	1:B:18:GLU:CB	2.61	0.46
1:B:512:TRP:C	1:B:514:VAL:H	2.24	0.46
1:A:12:LEU:HD23	1:A:66:MET:HB3	1.97	0.46
1:A:141:LYS:HG3	1:A:240:LYS:HE3	1.98	0.46
1:A:298:GLY:HA3	1:A:484:PHE:O	2.16	0.46
1:B:112:VAL:O	1:B:113:ASP:CB	2.39	0.46
1:A:112:VAL:O	1:A:113:ASP:CB	2.57	0.46
1:A:303:ASP:OD1	1:A:495:HIS:HE1	1.97	0.46
1:A:537:HIS:HB2	1:A:542:ASN:ND2	2.31	0.46
1:B:334:THR:HG22	1:B:377:ILE:HB	1.97	0.46
1:B:387:GLN:HA	1:B:391:TYR:CG	2.50	0.46
1:A:424:LEU:HD23	1:B:186:ILE:HG23	1.98	0.45
1:A:534:VAL:HG21	1:A:545:ILE:HG21	1.97	0.45
1:B:537:HIS:HB2	1:B:542:ASN:ND2	2.31	0.45
1:A:6:ASN:HA	1:A:7:PRO:HD3	1.72	0.45
1:A:144:THR:HG21	1:A:201:GLU:HB2	1.98	0.45
1:A:170:LEU:HD11	1:A:486:LEU:HD23	1.98	0.45
1:A:138:GLY:O	1:A:139:ASP:HB3	2.16	0.45
1:A:471:ARG:HA	1:A:472:PRO:HD2	1.85	0.45
1:B:26:ARG:HA	1:B:436:MET:CE	2.46	0.45
1:B:369:ARG:O	1:B:369:ARG:HD3	2.16	0.45
1:A:8:GLN:NE2	1:A:8:GLN:N	2.62	0.45
1:A:420:HIS:CD2	1:B:186:ILE:O	2.70	0.45
1:A:58:LEU:N	1:A:58:LEU:HD22	2.31	0.45
1:A:303:ASP:HA	1:A:495:HIS:HE1	1.81	0.45
1:A:113:ASP:OD1	3:A:724:HOH:O	2.11	0.45
1:B:186:ILE:HB	1:B:216:GLU:HG3	1.99	0.45
1:B:58:LEU:HD23	1:B:472:PRO:HB3	1.99	0.44
1:B:335:GLN:HB3	1:B:378:VAL:HG22	1.99	0.44
1:B:385:ASN:H	1:B:385:ASN:ND2	2.15	0.44
1:B:415:ILE:HD11	1:B:420:HIS:CG	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:462:LEU:HB3	1:B:463:PRO:HD3	1.99	0.44
1:A:103:ARG:HD3	1:A:299:ALA:HB3	1.99	0.44
1:A:141:LYS:HG3	1:A:240:LYS:CD	2.48	0.44
1:A:26:ARG:HA	1:A:436:MET:CE	2.47	0.44
1:A:434:ALA:HB1	1:B:516:LEU:HD13	2.00	0.44
1:B:415:ILE:HD12	1:B:415:ILE:O	2.17	0.44
1:A:103:ARG:HD3	1:A:299:ALA:HB1	1.99	0.44
1:B:17:ARG:CB	1:B:17:ARG:HH11	2.31	0.44
1:B:104:ASN:HD22	1:B:118:MET:HE2	1.83	0.44
1:B:117:VAL:HG22	1:B:120:GLU:OE1	2.18	0.44
1:A:144:THR:HB	1:A:201:GLU:CB	2.48	0.44
1:A:265:PHE:CZ	1:A:279:ILE:HD11	2.53	0.44
1:B:104:ASN:ND2	1:B:118:MET:HE2	2.33	0.44
1:A:44:ASN:HD22	1:A:44:ASN:C	2.25	0.44
1:A:379:TRP:CH2	1:B:356:MET:HE1	2.53	0.44
1:A:516:LEU:HD13	1:B:434:ALA:HB1	2.00	0.44
1:A:405:PHE:HB3	1:A:428:PHE:CE1	2.53	0.44
1:B:5:ARG:HH11	1:B:5:ARG:HB3	1.82	0.44
1:B:116:ASP:OD2	1:B:119:PRO:HD3	2.18	0.43
1:A:52:LEU:CD2	1:A:52:LEU:C	2.91	0.43
1:A:171:LYS:HD2	1:A:286:HIS:HE1	1.83	0.43
1:A:384:THR:O	1:A:387:GLN:HB2	2.18	0.43
1:B:92:THR:OG1	1:B:510:ASP:OD1	2.37	0.43
1:A:325:TRP:O	1:A:329:CYS:HB2	2.19	0.43
1:A:99:HIS:ND1	1:A:103:ARG:NH2	2.64	0.43
1:B:102:LEU:CD1	1:B:277:SER:HB3	2.46	0.43
1:B:111:VAL:HG13	1:B:115:LYS:N	2.33	0.43
1:A:480:LYS:HG3	3:A:601:HOH:O	2.19	0.43
1:B:23:LEU:HD22	1:B:23:LEU:H	1.83	0.43
1:B:106:SER:O	1:B:107:ASN:HB2	2.18	0.43
1:A:42:THR:HG22	1:A:51:LEU:HD13	2.01	0.43
1:A:383:GLY:CA	1:B:185:ASN:HD21	2.26	0.43
1:A:389:ALA:O	1:B:357:GLU:HB2	2.19	0.43
1:B:405:PHE:HB3	1:B:428:PHE:CE1	2.54	0.43
1:B:461:LEU:O	1:B:465:LYS:HG3	2.19	0.43
1:B:74:ARG:HG2	1:B:74:ARG:HH11	1.84	0.43
1:B:471:ARG:HA	1:B:472:PRO:HD2	1.85	0.43
1:A:12:LEU:CD1	1:A:329:CYS:SG	3.07	0.43
1:A:173:TYR:CZ	1:A:483:PRO:HD3	2.54	0.42
1:B:90:ASN:ND2	1:B:90:ASN:C	2.76	0.42
1:A:83:MET:HE2	1:A:97:VAL:CG1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:PHE:CE1	1:B:35:ARG:HD3	2.55	0.42
1:B:44:ASN:HD22	1:B:45:THR:N	2.18	0.42
1:B:104:ASN:HB3	1:B:118:MET:CE	2.50	0.42
1:B:552:ARG:CB	1:B:552:ARG:HH11	2.32	0.42
1:A:228:PHE:C	1:A:230:LEU:H	2.27	0.42
1:A:248:ASN:CB	1:A:251:LYS:HG3	2.49	0.42
1:A:125:LEU:HD23	1:A:128:MET:CE	2.49	0.42
1:A:435:LEU:HD12	1:A:435:LEU:HA	1.92	0.42
1:A:460:LYS:O	1:A:460:LYS:HG3	2.19	0.42
1:A:44:ASN:HD22	1:A:46:ASN:H	1.65	0.42
1:A:68:LEU:HD13	1:A:312:GLU:HG3	2.00	0.42
1:A:272:ARG:HG3	1:A:273:TYR:CE1	2.54	0.42
1:A:322:LEU:O	1:A:325:TRP:HB3	2.20	0.42
1:B:8:GLN:H	1:B:8:GLN:CD	2.27	0.42
1:B:257:ILE:CG2	1:B:258:ASP:N	2.83	0.42
1:A:45:THR:C	1:A:47:HIS:H	2.27	0.42
1:A:191:ILE:HG12	1:A:195:LEU:HD13	2.01	0.42
1:A:263:PHE:CD1	1:A:263:PHE:N	2.88	0.42
1:A:406:LEU:HD12	1:A:406:LEU:N	2.34	0.42
1:A:415:ILE:HD12	1:A:415:ILE:O	2.19	0.42
1:B:45:THR:C	1:B:47:HIS:H	2.28	0.42
1:B:361:LYS:HZ3	1:B:510:ASP:CB	2.28	0.42
1:B:406:LEU:N	1:B:406:LEU:HD12	2.35	0.42
1:B:446:LYS:CE	1:B:449:GLN:HE22	2.32	0.42
1:A:12:LEU:HD11	1:A:325:TRP:CE3	2.55	0.42
1:B:127:LYS:NZ	3:B:788:HOH:O	2.53	0.42
1:A:103:ARG:NH1	1:A:303:ASP:OD1	2.53	0.42
1:A:258:ASP:HA	1:A:259:PRO:HD3	1.87	0.42
1:A:370:VAL:HG13	1:A:372:HIS:CE1	2.54	0.42
1:A:245:LEU:HA	1:A:263:PHE:O	2.20	0.42
1:A:352:GLN:O	1:A:356:MET:HB2	2.20	0.42
1:A:399:LYS:HD3	1:B:373:GLN:OE1	2.19	0.42
1:B:104:ASN:HD22	1:B:118:MET:CE	2.32	0.41
1:A:420:HIS:HE1	1:B:223:THR:OG1	2.03	0.41
1:B:352:GLN:O	1:B:356:MET:HB2	2.20	0.41
1:A:444:ALA:O	1:A:448:LEU:HD23	2.20	0.41
1:A:41:LEU:CD1	1:A:43:LEU:HG	2.50	0.41
1:A:125:LEU:HA	1:A:125:LEU:HD23	1.92	0.41
1:A:407:ILE:HD11	1:A:424:LEU:HD12	2.03	0.41
1:B:249:THR:HG22	1:B:262:MET:SD	2.61	0.41
1:A:104:ASN:HB3	1:A:118:MET:CE	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:LYS:O	1:A:229:LEU:HB2	2.20	0.41
1:B:17:ARG:C	1:B:19:HIS:H	2.28	0.41
1:B:194:THR:O	1:B:198:LEU:HG	2.21	0.41
1:B:317:VAL:O	1:B:321:MET:HG2	2.21	0.41
1:B:322:LEU:O	1:B:325:TRP:HB3	2.20	0.41
1:A:173:TYR:HB3	1:A:291:ASN:ND2	2.35	0.41
1:A:186:ILE:HB	1:A:216:GLU:CG	2.51	0.41
1:A:520:LEU:HD12	1:B:434:ALA:HB2	2.02	0.41
1:B:186:ILE:HB	1:B:216:GLU:CG	2.51	0.41
1:B:234:ASP:HA	1:B:235:PRO:HD2	1.96	0.41
1:B:370:VAL:HG13	1:B:372:HIS:CE1	2.55	0.41
1:B:514:VAL:O	1:B:514:VAL:HG12	2.19	0.41
1:A:25:LEU:O	1:A:29:PHE:HD1	2.04	0.41
1:A:415:ILE:HD11	1:A:420:HIS:ND1	2.36	0.41
1:A:547:PHE:CZ	1:A:551:GLN:HG3	2.56	0.41
1:B:4:THR:HG23	1:B:372:HIS:HD2	1.86	0.41
1:B:8:GLN:CD	1:B:8:GLN:N	2.79	0.41
1:B:242:PHE:HB3	1:B:257:ILE:HD11	2.02	0.41
1:B:15:TRP:O	1:B:16:HIS:C	2.64	0.40
1:B:90:ASN:HD22	1:B:507:ASN:HD21	1.59	0.40
1:B:276:TRP:CZ3	1:B:299:ALA:HB2	2.56	0.40
1:A:351:PHE:HB3	3:A:746:HOH:O	2.21	0.40
1:A:74:ARG:NH2	1:A:505:ASP:HB2	2.37	0.40
1:A:248:ASN:HB3	1:A:251:LYS:CG	2.52	0.40
1:B:225:LYS:O	1:B:229:LEU:HB2	2.22	0.40
1:B:228:PHE:C	1:B:230:LEU:H	2.29	0.40
1:A:305:HIS:HE1	1:A:315:ALA:H	1.68	0.40
1:A:454:SER:O	1:A:455:PRO:C	2.65	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	553/557 (99%)	514 (93%)	34 (6%)	5 (1%)	14	27
1	B	553/557 (99%)	511 (92%)	35 (6%)	7 (1%)	9	18
All	All	1106/1114 (99%)	1025 (93%)	69 (6%)	12 (1%)	11	22

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	TYR
1	A	138	GLY
1	B	113	ASP
1	A	113	ASP
1	A	175	SER
1	B	16	HIS
1	B	22	GLU
1	B	2	ALA
1	B	3	LEU
1	B	383	GLY
1	A	200	PRO
1	B	200	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	475/478 (99%)	450 (95%)	25 (5%)	20	42
1	B	475/478 (99%)	448 (94%)	27 (6%)	18	39
All	All	950/956 (99%)	898 (94%)	52 (6%)	19	40

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	52	LEU
1	A	70	LEU
1	A	90	ASN

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Mol	Chain	Res	Type
1	A	115	LYS
1	A	117	VAL
1	A	123	LYS
1	A	133	GLN
1	A	141	LYS
1	A	161	LEU
1	A	229	LEU
1	A	233	LYS
1	A	243	VAL
1	A	254	GLU
1	A	265	PHE
1	A	296	LEU
1	A	344	LEU
1	A	346	ARG
1	A	411	THR
1	A	419	LEU
1	A	435	LEU
1	A	436	MET
1	A	456	GLU
1	A	514	VAL
1	A	516	LEU
1	B	8	GLN
1	B	52	LEU
1	B	56	LYS
1	B	70	LEU
1	B	74	ARG
1	B	81	GLU
1	B	90	ASN
1	B	105	ARG
1	B	115	LYS
1	B	117	VAL
1	B	143	TYR
1	B	144	THR
1	B	161	LEU
1	B	229	LEU
1	B	243	VAL
1	B	253	LYS
1	B	265	PHE
1	B	344	LEU
1	B	346	ARG
1	B	369	ARG
1	B	419	LEU

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Mol	Chain	Res	Type
1	B	435	LEU
1	B	436	MET
1	B	459	MET
1	B	471	ARG
1	B	516	LEU
1	B	552	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	13	GLN
1	A	44	ASN
1	A	90	ASN
1	A	215	GLN
1	A	261	ASN
1	A	286	HIS
1	A	352	GLN
1	A	385	ASN
1	A	387	GLN
1	A	395	HIS
1	A	410	GLN
1	A	413	HIS
1	A	420	HIS
1	A	495	HIS
1	A	511	GLN
1	A	542	ASN
1	A	550	GLN
1	B	6	ASN
1	B	8	GLN
1	B	10	GLN
1	B	16	HIS
1	B	24	ASN
1	B	27	HIS
1	B	44	ASN
1	B	49	HIS
1	B	90	ASN
1	B	99	HIS
1	B	104	ASN
1	B	107	ASN
1	B	122	ASN
1	B	215	GLN

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Mol	Chain	Res	Type
1	B	294	GLN
1	B	352	GLN
1	B	353	GLN
1	B	359	ASN
1	B	385	ASN
1	B	387	GLN
1	B	395	HIS
1	B	410	GLN
1	B	412	GLN
1	B	413	HIS
1	B	420	HIS
1	B	449	GLN
1	B	495	HIS
1	B	519	GLN
1	B	542	ASN
1	B	550	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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	6PG	B	601	-	16,16,16	1.42	3 (18%)	21,23,23	0.93	1 (4%)
2	6PG	A	600	-	16,16,16	1.36	4 (25%)	21,23,23	0.97	1 (4%)

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	6PG	B	601	-	-	9/22/22/22	-
2	6PG	A	600	-	-	14/22/22/22	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	6PG	O1-C1	-3.37	1.19	1.30
2	A	600	6PG	O1-C1	-2.90	1.21	1.30
2	B	601	6PG	C3-C2	2.37	1.58	1.53
2	A	600	6PG	P-O6	-2.34	1.52	1.60
2	A	600	6PG	C4-C3	2.15	1.57	1.53
2	A	600	6PG	P-O1P	2.09	1.57	1.50
2	B	601	6PG	P-O6	-2.03	1.53	1.60

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	6PG	O1-C1-C2	2.33	119.80	113.31
2	A	600	6PG	O1-C1-C2	2.24	119.53	113.31

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	600	6PG	C4-C5-C6-O6
2	A	600	6PG	O5-C5-C6-O6
2	B	601	6PG	C1-C2-C3-C4
2	B	601	6PG	C1-C2-C3-O3
2	B	601	6PG	O2-C2-C3-C4
2	B	601	6PG	O2-C2-C3-O3
2	B	601	6PG	C2-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
2	A	600	6PG	O4-C4-C5-C6
2	A	600	6PG	C3-C4-C5-C6
2	B	601	6PG	C2-C3-C4-O4
2	B	601	6PG	O3-C3-C4-O4
2	B	601	6PG	O3-C3-C4-C5
2	A	600	6PG	C3-C4-C5-O5
2	A	600	6PG	O4-C4-C5-O5
2	A	600	6PG	O1-C1-C2-O2
2	B	601	6PG	C4-C5-C6-O6
2	A	600	6PG	C2-C3-C4-O4
2	A	600	6PG	O1A-C1-C2-C3
2	A	600	6PG	O1-C1-C2-C3
2	A	600	6PG	O1A-C1-C2-O2
2	A	600	6PG	O3-C3-C4-O4
2	A	600	6PG	C2-C3-C4-C5
2	A	600	6PG	O3-C3-C4-C5

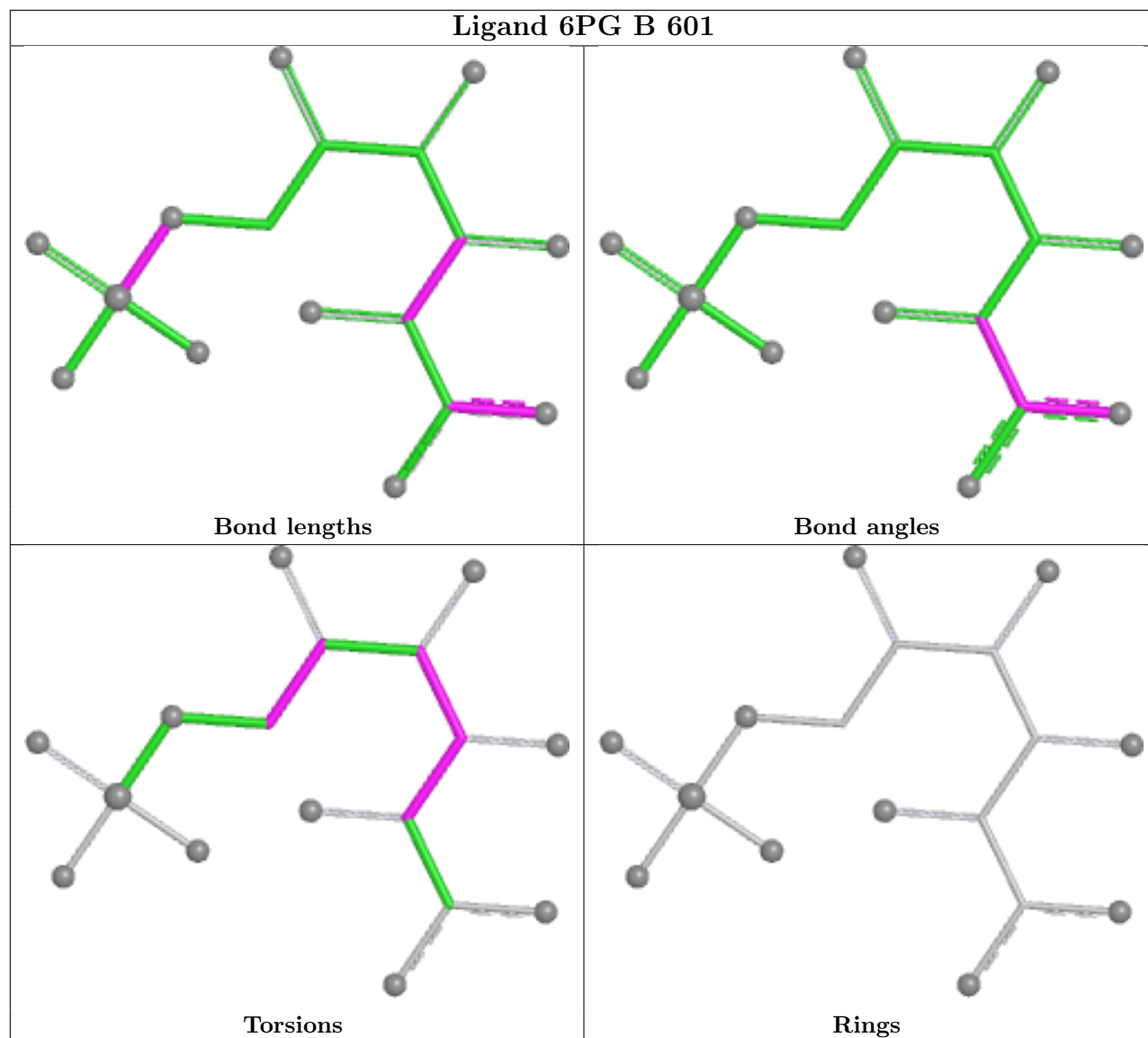
There are no ring outliers.

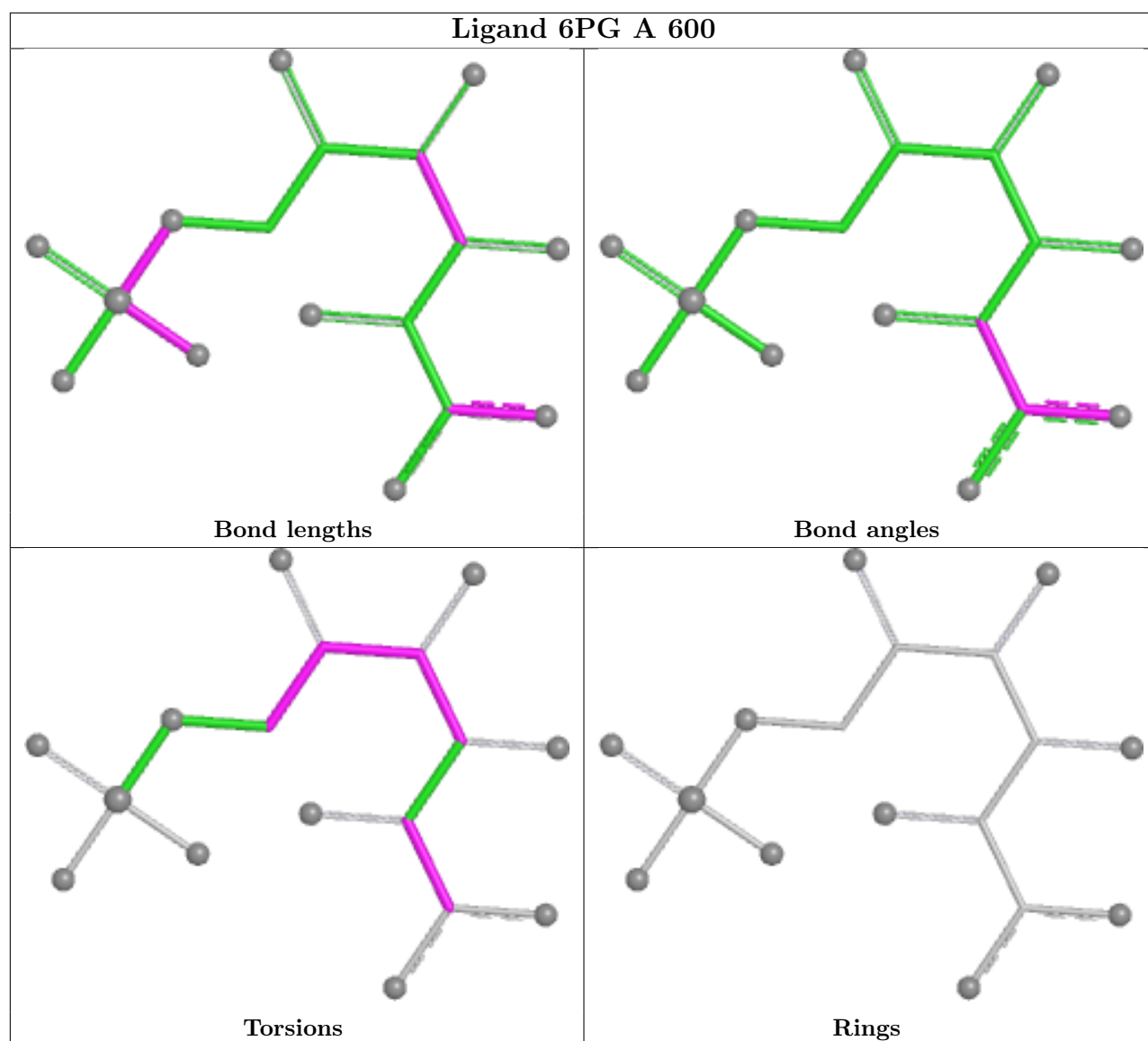
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	6PG	1	0
2	A	600	6PG	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.

Ligand 6PG B 601





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	555/557 (99%)	-0.03	17 (3%) 51 47	10, 27, 61, 87	0
1	B	555/557 (99%)	-0.14	13 (2%) 61 57	9, 25, 52, 86	0
All	All	1110/1114 (99%)	-0.08	30 (2%) 56 51	9, 26, 57, 87	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	175	SER	4.9
1	A	137	SER	4.4
1	B	18	GLU	4.3
1	A	138	GLY	3.8
1	B	20	GLY	3.4
1	A	232	ALA	3.3
1	B	19	HIS	3.3
1	A	452	GLY	3.2
1	B	17	ARG	2.9
1	A	175	SER	2.9
1	A	30	ASP	2.8
1	B	383	GLY	2.7
1	B	143	TYR	2.7
1	A	249	THR	2.5
1	A	112	VAL	2.5
1	A	454	SER	2.5
1	B	112	VAL	2.5
1	A	455	PRO	2.3
1	B	21	SER	2.3
1	A	176	GLY	2.3
1	A	510	ASP	2.3
1	A	383	GLY	2.3
1	A	144	THR	2.3
1	A	449	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	254	GLU	2.2
1	B	113	ASP	2.2
1	B	253	LYS	2.2
1	B	232	ALA	2.0
1	A	174	SER	2.0
1	A	451	ALA	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](#)

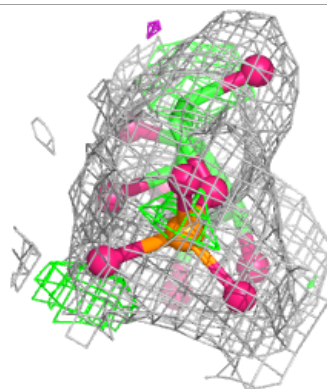
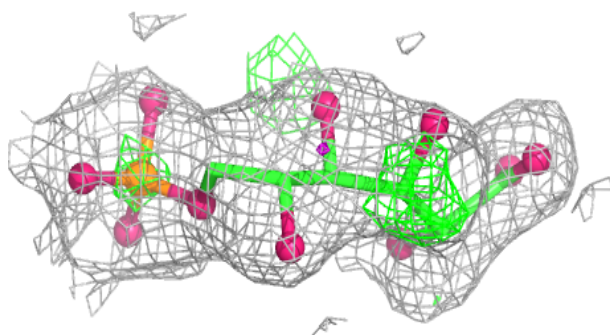
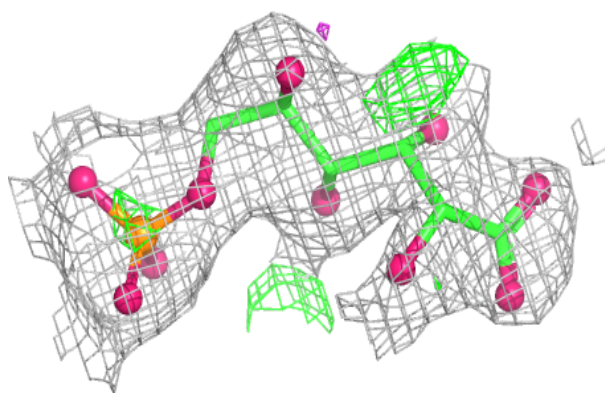
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	6PG	B	601	17/17	0.91	0.11	31,41,49,54	0
2	6PG	A	600	17/17	0.92	0.19	42,57,73,74	0

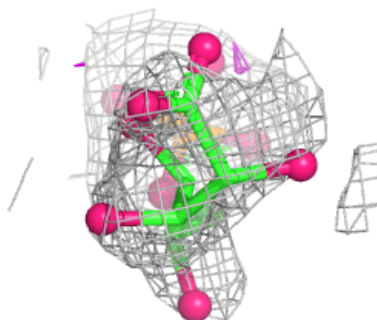
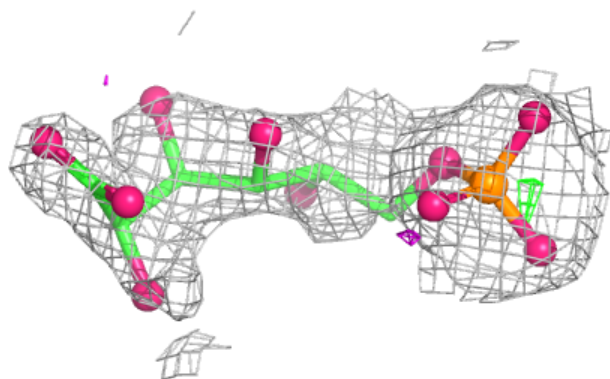
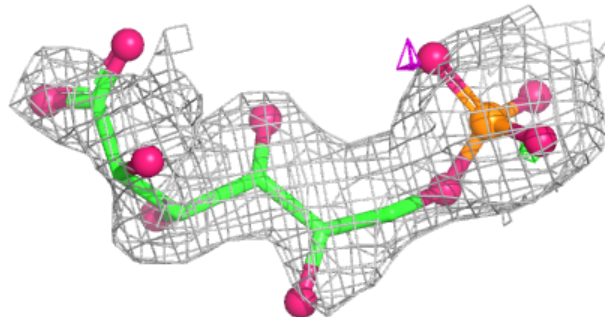
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6PG B 601:

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

**Electron density around 6PG A 600:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.