



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 05:37 AM UTC

PDB ID : 6JJ7 / pdb_00006jj7
Title : Crystal structure of OsHXK6-Glc complex
Authors : He, C.; Wei, P.; Chen, J.; Wang, H.; Wan, Y.; Zhou, J.; Zhu, Y.; Huang, W.; Yin, L.
Deposited on : 2019-02-25
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

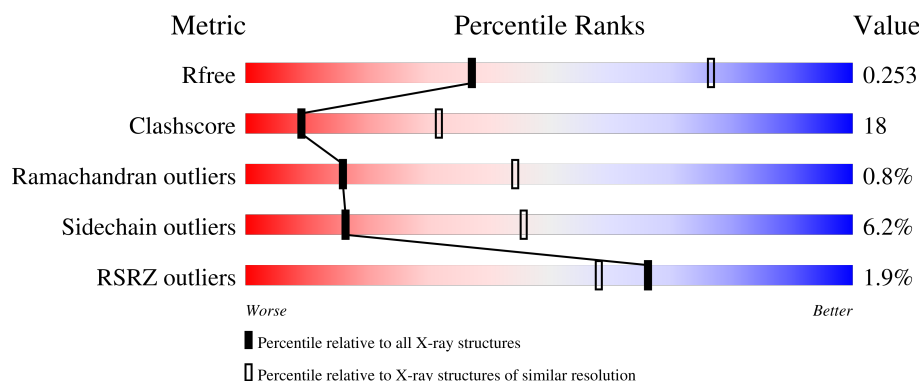
1 Overall quality at a glance

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	465	2% 63% 30% 5% •
1	C	465	2% 66% 28% 5% •
1	E	465	2% 70% 26% • •

2 Entry composition [i](#)

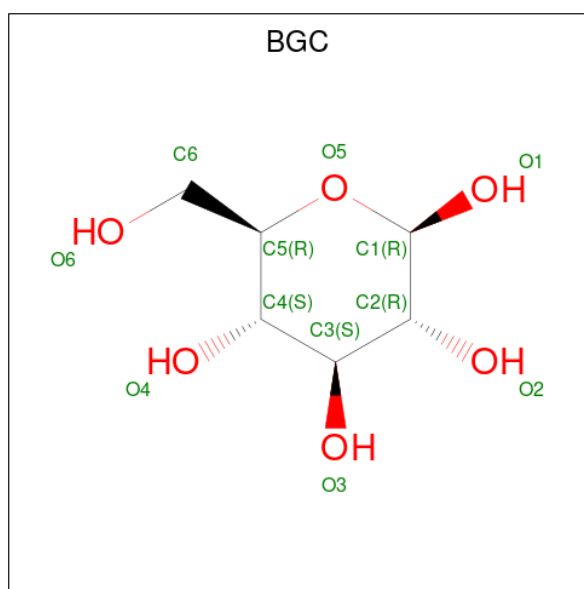
There are 2 unique types of molecules in this entry. The entry contains 10803 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Rice hexokinase 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3589	2257	634	683	15			
1	C	465	Total	C	N	O	S	0	0	0
			3589	2257	634	683	15			
1	E	465	Total	C	N	O	S	0	0	0
			3589	2257	634	683	15			

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆) (labeled as "Ligand of Interest" by depositor).

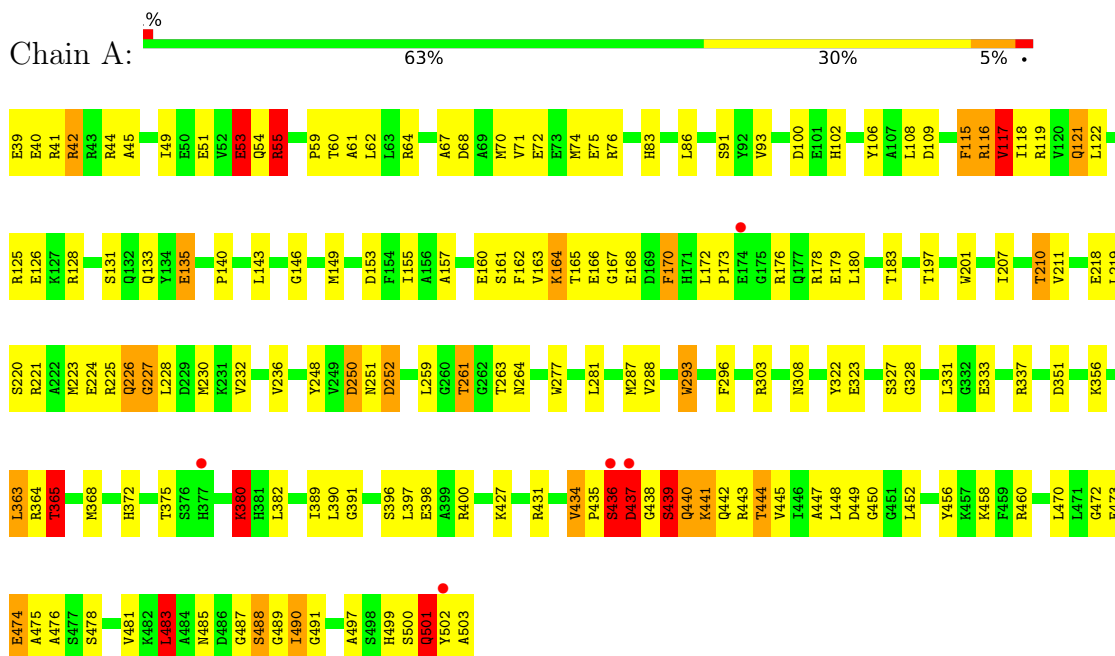


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	C	1	Total	C	O	0	0
			12	6	6		
2	E	1	Total	C	O	0	0
			12	6	6		

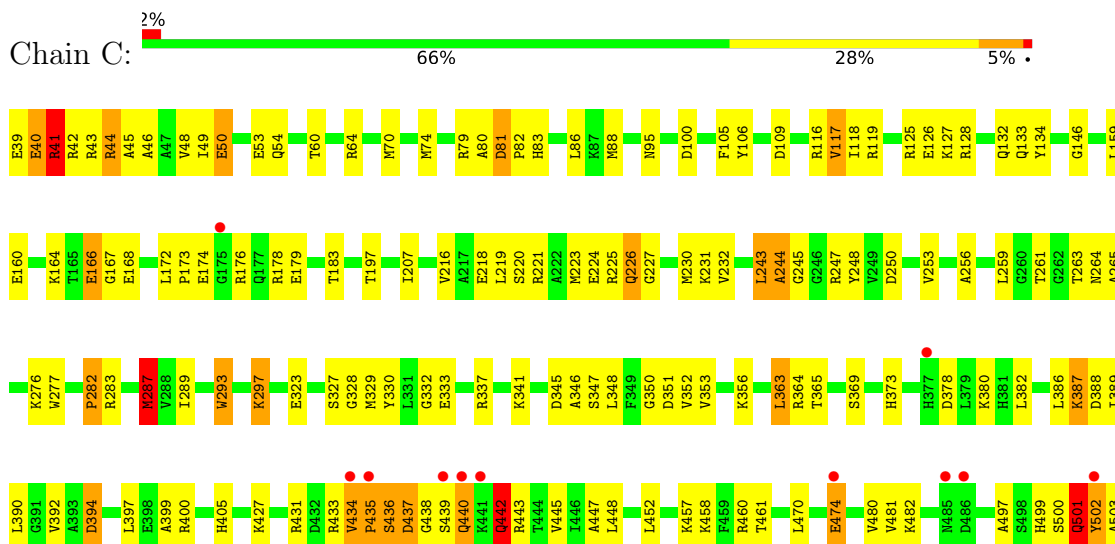
3 Residue-property plots

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.

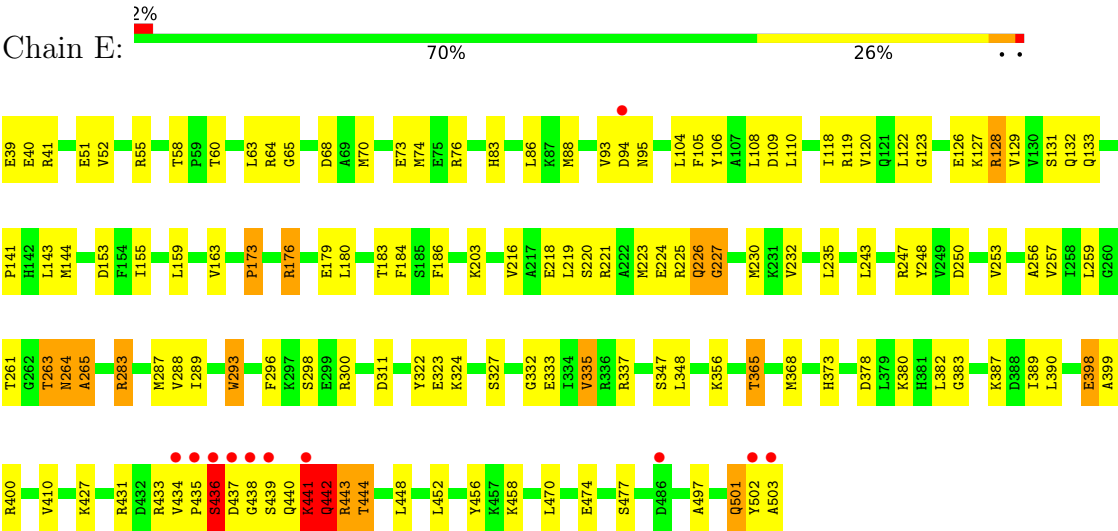
• Molecule 1: Rice hexokinase 6



• Molecule 1: Rice hexokinase 6



• Molecule 1: Rice hexokinase 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.57Å 131.57Å 188.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.47 – 2.90 45.47 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.47-2.90) 97.0 (45.47-2.90)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.15 (at 2.91Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.200 , 0.248 0.210 , 0.253	Depositor DCC
R_{free} test set	1997 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	39.8	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 21.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10803	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% 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 [i](#)

5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.19	19/3652 (0.5%)	1.01	17/4936 (0.3%)
1	C	0.99	6/3652 (0.2%)	0.98	19/4936 (0.4%)
1	E	0.93	9/3652 (0.2%)	0.91	10/4936 (0.2%)
All	All	1.04	34/10956 (0.3%)	0.97	46/14808 (0.3%)

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	MET	C-O	-6.42	1.16	1.24
1	C	287	MET	C-O	-6.38	1.16	1.23
1	A	250	ASP	C-O	-6.03	1.16	1.24
1	A	363	LEU	C-O	-5.91	1.17	1.24
1	E	263	THR	C-O	-5.82	1.17	1.23
1	A	135	GLU	C-O	-5.78	1.17	1.23
1	E	442	GLN	C-O	-5.61	1.16	1.23
1	A	460	ARG	C-O	-5.56	1.17	1.24
1	E	203	LYS	CA-C	-5.49	1.48	1.53
1	E	443	ARG	C-O	-5.43	1.17	1.24
1	A	72	GLU	C-O	-5.40	1.17	1.24
1	E	55	ARG	C-O	-5.40	1.17	1.24
1	A	157	ALA	C-O	-5.40	1.17	1.24
1	A	476	ALA	C-O	-5.34	1.17	1.24
1	C	351	ASP	CA-C	-5.31	1.45	1.52
1	E	203	LYS	C-O	-5.26	1.19	1.24
1	A	364	ARG	C-O	-5.25	1.17	1.23
1	A	250	ASP	CA-C	-5.24	1.46	1.52
1	C	364	ARG	C-O	-5.24	1.17	1.23
1	A	116	ARG	C-O	-5.22	1.17	1.23
1	A	170	PHE	CA-C	-5.19	1.46	1.52
1	C	363	LEU	C-O	-5.18	1.17	1.24
1	E	264	ASN	C-O	-5.13	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	53	GLU	C-O	-5.12	1.18	1.24
1	A	117	VAL	C-O	-5.11	1.18	1.24
1	A	115	PHE	C-O	-5.10	1.17	1.23
1	C	81	ASP	CA-C	-5.08	1.47	1.53
1	E	265	ALA	C-O	-5.07	1.17	1.23
1	A	491	GLY	C-O	-5.07	1.17	1.23
1	C	353	VAL	N-CA	-5.05	1.41	1.47
1	A	121	GLN	C-O	-5.04	1.18	1.24
1	E	52	VAL	C-O	-5.04	1.18	1.24
1	A	76	ARG	C-O	-5.01	1.18	1.24
1	A	55	ARG	C-O	-5.01	1.17	1.24

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	127	LYS	N-CA-C	13.93	126.15	110.97
1	A	437	ASP	N-CA-C	-10.10	96.29	111.34
1	C	474	GLU	CA-CB-CG	9.90	133.90	114.10
1	A	227	GLY	N-CA-C	-9.47	85.83	113.30
1	A	226	GLN	N-CA-C	8.96	124.34	113.41
1	E	52	VAL	N-CA-C	-8.71	102.34	110.53
1	C	502	TYR	N-CA-C	-8.64	86.82	111.00
1	C	501	GLN	N-CA-C	8.08	124.13	113.30
1	C	167	GLY	N-CA-C	-7.90	90.39	113.30
1	A	210	THR	CB-CA-C	7.74	122.39	109.38
1	C	226	GLN	N-CA-C	7.55	122.79	113.50
1	A	488	SER	N-CA-C	7.42	120.25	111.71
1	E	226	GLN	N-CA-C	7.39	122.58	113.50
1	A	490	ILE	N-CA-C	-7.21	103.44	110.72
1	C	474	GLU	CB-CG-CD	7.04	124.56	112.60
1	A	41	ARG	N-CA-C	-7.00	103.58	111.07
1	C	48	VAL	N-CA-C	-6.94	105.00	111.45
1	E	129	VAL	N-CA-C	-6.75	97.46	107.51
1	A	380	LYS	N-CA-C	6.52	118.39	111.28
1	C	174	GLU	CA-CB-CG	-6.48	101.14	114.10
1	A	501	GLN	N-CA-C	-6.43	105.73	112.93
1	C	282	PRO	N-CA-C	6.41	122.39	111.68
1	E	501	GLN	N-CA-C	6.25	120.94	113.12
1	C	117	VAL	N-CA-C	-6.11	98.83	107.75
1	C	297	LYS	N-CA-C	-6.03	98.78	108.55
1	A	483	LEU	N-CA-C	-5.95	100.30	109.76
1	C	482	LYS	N-CA-C	5.94	117.43	108.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	166	GLU	N-CA-C	-5.94	99.60	109.46
1	A	434	VAL	CA-C-N	-5.85	113.91	119.76
1	A	434	VAL	C-N-CA	-5.85	113.91	119.76
1	A	436	SER	N-CA-C	-5.77	94.83	111.00
1	E	227	GLY	N-CA-C	-5.77	96.57	113.30
1	C	40	GLU	N-CA-C	-5.63	106.25	113.01
1	C	134	TYR	N-CA-C	5.50	117.55	109.24
1	C	244	ALA	N-CA-C	-5.40	103.30	111.56
1	C	501	GLN	CA-C-N	5.39	131.40	121.70
1	C	501	GLN	C-N-CA	5.39	131.40	121.70
1	E	127	LYS	CA-C-N	5.39	131.40	121.70
1	E	127	LYS	C-N-CA	5.39	131.40	121.70
1	A	365	THR	CB-CA-C	5.31	120.63	110.17
1	A	167	GLY	N-CA-C	-5.28	97.99	113.30
1	C	442	GLN	CA-CB-CG	5.26	124.62	114.10
1	A	61	ALA	N-CA-C	-5.25	105.64	111.36
1	E	398	GLU	N-CA-C	-5.22	105.27	111.69
1	A	54	GLN	N-CA-C	5.20	117.86	111.82
1	E	126	GLU	N-CA-C	5.01	119.52	113.41

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	3589	0	3582	127	0
1	C	3589	0	3582	136	0
1	E	3589	0	3582	134	0
2	A	12	0	11	0	0
2	C	12	0	11	0	0
2	E	12	0	11	2	0
All	All	10803	0	10779	393	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 18.

All (393) 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:C:437:ASP:HB2	1:C:438:GLY:CA	1.14	1.49
1:E:440:GLN:HA	1:E:441:LYS:CB	1.55	1.33
1:E:176:ARG:NH2	1:E:502:TYR:O	1.63	1.27
1:A:440:GLN:HE21	1:A:441:LYS:N	1.31	1.27
1:E:440:GLN:CA	1:E:441:LYS:HB2	1.50	1.25
1:E:437:ASP:OD1	1:E:439:SER:CB	1.86	1.22
1:C:437:ASP:CB	1:C:438:GLY:CA	2.02	1.22
1:C:437:ASP:CB	1:C:438:GLY:HA2	1.64	1.20
1:E:437:ASP:OD1	1:E:439:SER:HB3	1.39	1.19
1:A:252:ASP:OD2	1:A:443:ARG:O	1.62	1.14
1:A:119:ARG:NH2	1:A:133:GLN:OE1	1.82	1.12
1:A:438:GLY:H	1:A:439:SER:HB2	1.11	1.09
1:C:437:ASP:HB2	1:C:438:GLY:HA3	1.26	1.09
1:A:39:GLU:OE2	1:A:40:GLU:N	1.87	1.06
1:C:119:ARG:NH2	1:C:133:GLN:OE1	1.88	1.05
1:C:39:GLU:HG3	1:C:41:ARG:NH2	1.75	1.01
1:A:438:GLY:N	1:A:439:SER:HB2	1.76	1.01
1:C:437:ASP:CB	1:C:438:GLY:HA3	1.80	0.99
1:A:440:GLN:NE2	1:A:441:LYS:H	1.61	0.98
1:A:435:PRO:HA	1:A:436:SER:HB2	1.48	0.95
1:A:160:GLU:OE1	1:A:164:LYS:NZ	1.99	0.95
1:A:440:GLN:HE21	1:A:441:LYS:H	0.98	0.94
1:E:442:GLN:H	1:E:442:GLN:HE21	1.07	0.94
1:E:501:GLN:HB2	1:E:503:ALA:H	1.31	0.93
1:A:440:GLN:NE2	1:A:441:LYS:N	2.14	0.91
1:A:328:GLY:O	1:A:365:THR:CG2	2.19	0.90
1:C:106:TYR:OH	1:C:166:GLU:OE1	1.90	0.89
1:C:166:GLU:OE2	1:C:178:ARG:NH1	2.06	0.89
1:C:44:ARG:HG2	1:C:44:ARG:HH21	1.35	0.88
1:A:435:PRO:CA	1:A:436:SER:HB2	2.03	0.88
1:A:328:GLY:O	1:A:365:THR:HG21	1.73	0.88
1:C:45:ALA:O	1:C:49:ILE:HD12	1.74	0.87
1:A:117:VAL:HG23	1:A:135:GLU:HB3	1.58	0.86
1:E:119:ARG:NH2	1:E:133:GLN:OE1	2.09	0.84
1:E:436:SER:H	1:E:437:ASP:C	1.85	0.84
1:E:442:GLN:H	1:E:442:GLN:NE2	1.77	0.83
1:A:179:GLU:OE2	1:A:501:GLN:HG2	1.78	0.83
1:C:39:GLU:HG3	1:C:41:ARG:HH21	1.40	0.82
1:A:121:GLN:OE1	1:A:170:PHE:HA	1.79	0.81
1:A:389:ILE:HG22	1:A:390:LEU:HD23	1.60	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:434:VAL:HG13	1:C:435:PRO:HD2	1.61	0.80
1:A:133:GLN:HG3	1:A:170:PHE:HZ	1.47	0.78
1:C:39:GLU:CG	1:C:41:ARG:NH2	2.45	0.78
1:E:436:SER:N	1:E:437:ASP:HB3	1.98	0.78
1:C:49:ILE:O	1:C:53:GLU:HG3	1.84	0.78
1:C:431:ARG:NH2	1:C:442:GLN:HE21	1.81	0.78
1:E:437:ASP:OD1	1:E:439:SER:N	2.16	0.78
1:A:100:ASP:OD1	1:A:125:ARG:HB2	1.83	0.77
1:A:440:GLN:HE21	1:A:440:GLN:C	1.93	0.77
1:A:438:GLY:CA	1:A:439:SER:HB2	2.15	0.77
1:C:172:LEU:HD11	1:C:178:ARG:HG2	1.67	0.77
1:C:332:GLY:H	1:C:365:THR:HG22	1.50	0.77
1:C:100:ASP:OD1	1:C:125:ARG:HB3	1.84	0.77
1:C:60:THR:O	1:C:64:ARG:HG3	1.84	0.77
1:C:159:LEU:HD23	1:C:223:MET:HE3	1.65	0.76
1:C:387:LYS:HE2	1:C:394:ASP:HA	1.68	0.76
1:A:39:GLU:CD	1:A:40:GLU:H	1.92	0.76
1:C:44:ARG:HG2	1:C:44:ARG:NH2	1.97	0.76
1:A:128:ARG:NH1	1:A:248:TYR:O	2.19	0.75
1:E:437:ASP:OD1	1:E:439:SER:HB2	1.82	0.75
1:E:437:ASP:CG	1:E:439:SER:N	2.44	0.75
1:C:261:THR:O	1:C:327:SER:OG	2.03	0.75
1:A:180:LEU:HD23	1:A:232:VAL:HG22	1.68	0.75
1:C:40:GLU:O	1:C:40:GLU:HG2	1.87	0.74
1:A:287:MET:HE3	1:A:288:VAL:H	1.52	0.74
1:E:438:GLY:H	1:E:441:LYS:NZ	1.86	0.73
1:E:261:THR:O	1:E:327:SER:OG	2.06	0.73
1:C:216:VAL:HG13	1:C:232:VAL:HG11	1.71	0.73
1:C:277:TRP:CE2	1:C:287:MET:HE1	2.23	0.73
1:C:382:LEU:HD23	1:C:382:LEU:O	1.90	0.72
1:C:243:LEU:HD12	1:C:243:LEU:O	1.89	0.71
1:E:93:VAL:N	1:E:287:MET:HE1	2.05	0.71
1:E:442:GLN:HE21	1:E:442:GLN:N	1.87	0.71
1:E:332:GLY:H	1:E:365:THR:HG22	1.56	0.71
1:E:64:ARG:NH2	1:E:470:LEU:O	2.23	0.71
1:C:437:ASP:HB2	1:C:438:GLY:HA2	0.72	0.71
1:E:163:VAL:HG11	1:E:230:MET:HE1	1.74	0.70
1:E:287:MET:HE3	1:E:288:VAL:H	1.57	0.70
1:E:173:PRO:HB2	1:E:176:ARG:HE	1.57	0.70
1:E:176:ARG:HH22	1:E:502:TYR:CB	2.05	0.69
1:A:448:LEU:HD13	1:A:481:VAL:HG13	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:GLN:HG3	1:A:170:PHE:CZ	2.28	0.69
1:A:207:ILE:O	1:A:210:THR:HG22	1.93	0.69
1:E:440:GLN:HG2	1:E:441:LYS:HB3	1.74	0.69
1:C:277:TRP:CD2	1:C:287:MET:HE1	2.27	0.68
1:A:51:GLU:OE1	1:A:55:ARG:NH1	2.26	0.68
1:E:263:THR:HG22	1:E:323:GLU:HG3	1.76	0.68
1:A:368:MET:HE3	1:A:456:TYR:HE2	1.59	0.68
1:A:93:VAL:N	1:A:287:MET:HE1	2.09	0.68
1:A:442:GLN:OE1	1:A:442:GLN:N	2.19	0.67
1:A:179:GLU:OE1	1:A:500:SER:OG	2.10	0.67
1:C:434:VAL:HG13	1:C:435:PRO:CD	2.24	0.67
1:E:368:MET:HE3	1:E:456:TYR:HE2	1.58	0.66
1:C:218:GLU:OE2	1:C:221:ARG:NH2	2.27	0.66
1:C:81:ASP:OD2	1:C:82:PRO:HD2	1.96	0.65
1:C:373:HIS:HE1	1:C:458:LYS:H	1.43	0.65
1:C:64:ARG:NH2	1:C:470:LEU:O	2.29	0.65
1:A:64:ARG:NH2	1:A:470:LEU:O	2.28	0.65
1:A:372:HIS:O	1:A:458:LYS:NZ	2.29	0.65
1:A:287:MET:HE3	1:A:288:VAL:N	2.12	0.65
1:E:437:ASP:OD1	1:E:439:SER:CA	2.43	0.65
1:C:128:ARG:NH1	1:C:248:TYR:O	2.29	0.64
1:E:60:THR:O	1:E:64:ARG:HG3	1.97	0.64
1:C:40:GLU:O	1:C:43:ARG:HB3	1.98	0.64
1:C:173:PRO:HG2	1:C:176:ARG:HH11	1.63	0.64
1:C:373:HIS:CE1	1:C:458:LYS:H	2.15	0.64
1:E:159:LEU:HD23	1:E:223:MET:HE3	1.80	0.64
1:A:45:ALA:O	1:A:49:ILE:HD12	1.97	0.63
1:E:259:LEU:O	1:E:452:LEU:HB2	1.98	0.63
1:A:431:ARG:HG2	1:A:444:THR:HG21	1.80	0.63
1:A:437:ASP:OD2	1:A:439:SER:OG	2.12	0.63
1:E:176:ARG:HH22	1:E:502:TYR:HB2	1.63	0.63
1:C:39:GLU:HG3	1:C:41:ARG:HH22	1.64	0.62
1:A:333:GLU:O	1:A:337:ARG:HG3	1.99	0.62
1:E:365:THR:HB	1:E:368:MET:HE2	1.81	0.61
1:A:220:SER:O	1:A:224:GLU:HG3	2.01	0.61
1:E:434:VAL:HG23	1:E:434:VAL:O	2.01	0.61
1:E:438:GLY:H	1:E:441:LYS:HZ1	1.48	0.61
1:A:42:ARG:NH2	1:A:398:GLU:OE1	2.34	0.61
1:A:49:ILE:O	1:A:53:GLU:HG3	2.01	0.61
1:E:431:ARG:HD2	1:E:444:THR:CG2	2.31	0.60
1:E:180:LEU:HD23	1:E:232:VAL:HG22	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:93:VAL:H	1:E:287:MET:HE1	1.62	0.60
1:E:94:ASP:OD1	1:E:95:ASN:N	2.31	0.60
1:E:436:SER:N	1:E:437:ASP:O	2.34	0.60
1:C:378:ASP:O	1:C:400:ARG:NH2	2.35	0.60
1:A:86:LEU:HD22	1:A:293:TRP:HA	1.84	0.60
1:A:162:PHE:O	1:A:165:THR:OG1	2.20	0.60
1:A:263:THR:HG22	1:A:323:GLU:HG3	1.81	0.60
1:E:68:ASP:OD1	1:E:427:LYS:NZ	2.30	0.60
1:C:382:LEU:HD23	1:C:382:LEU:C	2.27	0.59
1:C:435:PRO:C	1:C:437:ASP:H	2.10	0.59
1:E:226:GLN:N	1:E:227:GLY:HA2	2.16	0.59
1:E:250:ASP:OD2	1:E:443:ARG:NH1	2.36	0.59
1:E:501:GLN:HB2	1:E:503:ALA:N	2.10	0.59
1:A:223:MET:O	1:A:227:GLY:HA2	2.03	0.59
1:E:311:ASP:OD2	1:E:324:LYS:NZ	2.31	0.59
1:A:218:GLU:OE2	1:A:221:ARG:NH2	2.35	0.59
1:A:70:MET:O	1:A:74:MET:HG3	2.03	0.58
1:A:435:PRO:HA	1:A:436:SER:CB	2.28	0.58
1:C:176:ARG:NH2	1:C:502:TYR:O	2.36	0.58
1:E:141:PRO:HA	1:E:144:MET:HE3	1.84	0.58
1:C:39:GLU:CD	1:C:41:ARG:HH22	2.12	0.58
1:C:176:ARG:HD2	1:C:503:ALA:HA	1.85	0.58
1:E:438:GLY:N	1:E:441:LYS:NZ	2.52	0.58
1:C:44:ARG:HH21	1:C:44:ARG:CG	2.12	0.57
1:C:226:GLN:N	1:C:227:GLY:HA2	2.19	0.57
1:A:93:VAL:H	1:A:287:MET:HE1	1.69	0.57
1:E:128:ARG:HB2	1:E:248:TYR:OH	2.05	0.57
1:C:332:GLY:N	1:C:365:THR:HG22	2.19	0.56
1:A:153:ASP:OD1	1:A:225:ARG:NH1	2.38	0.56
1:A:502:TYR:O	1:A:503:ALA:HB2	2.06	0.56
1:C:259:LEU:O	1:C:452:LEU:HB2	2.05	0.56
1:C:41:ARG:O	1:C:43:ARG:N	2.38	0.56
1:C:347:SER:HB3	1:C:350:GLY:O	2.05	0.56
1:C:356:LYS:HE3	1:C:389:ILE:O	2.05	0.56
1:A:176:ARG:HD3	1:A:503:ALA:HA	1.88	0.55
1:A:382:LEU:HD22	1:A:400:ARG:HB3	1.87	0.55
1:E:356:LYS:HE3	1:E:389:ILE:O	2.06	0.55
1:E:220:SER:O	1:E:224:GLU:HG3	2.06	0.55
1:E:348:LEU:HD11	1:E:399:ALA:HA	1.87	0.55
1:E:436:SER:H	1:E:437:ASP:CA	2.18	0.55
1:A:179:GLU:OE2	1:A:501:GLN:CG	2.51	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:333:GLU:O	1:C:337:ARG:HG3	2.07	0.55
1:E:332:GLY:H	1:E:365:THR:CG2	2.20	0.55
1:C:431:ARG:NH2	1:C:442:GLN:NE2	2.52	0.54
1:E:382:LEU:HD22	1:E:400:ARG:HB3	1.90	0.54
1:A:146:GLY:O	1:A:207:ILE:HG23	2.07	0.54
1:C:176:ARG:CD	1:C:503:ALA:HA	2.37	0.54
1:E:436:SER:N	1:E:437:ASP:CB	2.70	0.54
1:A:356:LYS:HE3	1:A:389:ILE:O	2.06	0.54
1:C:126:GLU:HG2	1:C:127:LYS:HG3	1.88	0.54
1:C:282:PRO:O	1:C:283:ARG:C	2.48	0.54
1:E:435:PRO:HA	1:E:436:SER:CB	2.37	0.54
1:E:433:ARG:C	1:E:434:VAL:HG13	2.33	0.54
1:A:328:GLY:O	1:A:365:THR:HG23	2.06	0.53
1:A:434:VAL:CG2	1:A:441:LYS:HG2	2.39	0.53
1:A:331:LEU:HB3	1:A:368:MET:HE1	1.90	0.53
1:C:95:ASN:OD1	1:C:95:ASN:N	2.38	0.53
1:C:39:GLU:CG	1:C:41:ARG:HH22	2.18	0.53
1:C:88:MET:HE3	1:C:289:ILE:HG21	1.90	0.53
1:C:250:ASP:OD1	1:C:443:ARG:NH1	2.41	0.53
1:E:437:ASP:CG	1:E:439:SER:H	2.15	0.53
1:A:211:VAL:HG23	1:A:211:VAL:O	2.07	0.53
1:A:259:LEU:O	1:A:452:LEU:HB2	2.09	0.53
1:A:435:PRO:CB	1:A:436:SER:HB2	2.38	0.53
1:A:438:GLY:CA	1:A:439:SER:CB	2.84	0.52
1:A:40:GLU:O	1:A:40:GLU:HG2	2.09	0.52
1:A:449:ASP:OD1	1:A:450:GLY:N	2.41	0.52
1:A:437:ASP:CG	1:A:438:GLY:H	2.17	0.52
1:C:369:SER:O	1:C:373:HIS:HD2	1.92	0.52
1:E:106:TYR:CE2	1:E:163:VAL:HG12	2.45	0.52
1:A:440:GLN:NE2	1:A:440:GLN:CA	2.72	0.52
1:C:457:LYS:O	1:C:461:THR:HG23	2.10	0.52
1:A:108:LEU:HD21	1:A:155:ILE:HG21	1.92	0.52
1:C:106:TYR:CE2	1:C:119:ARG:HD3	2.45	0.52
1:C:79:ARG:O	1:C:80:ALA:HB3	2.09	0.52
1:C:243:LEU:HB2	1:C:256:ALA:HB2	1.92	0.51
1:C:178:ARG:HB2	1:C:230:MET:HE3	1.93	0.51
1:E:95:ASN:HD21	1:E:283:ARG:HD3	1.74	0.51
1:A:207:ILE:O	1:A:210:THR:CG2	2.58	0.51
1:A:440:GLN:NE2	1:A:440:GLN:HA	2.26	0.51
1:C:53:GLU:OE1	1:C:405:HIS:NE2	2.36	0.51
1:C:226:GLN:H	1:C:227:GLY:HA2	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:110:LEU:HD13	1:E:155:ILE:HD11	1.92	0.51
1:C:348:LEU:HD11	1:C:399:ALA:HA	1.91	0.51
1:E:440:GLN:HG2	1:E:441:LYS:CB	2.40	0.51
1:C:373:HIS:HE1	1:C:458:LYS:N	2.07	0.51
1:A:497:ALA:C	1:A:499:HIS:H	2.18	0.51
1:A:125:ARG:HG2	1:A:125:ARG:O	2.10	0.50
1:E:373:HIS:NE2	1:E:458:LYS:HG3	2.25	0.50
1:A:250:ASP:OD2	1:A:443:ARG:NH1	2.44	0.50
1:A:331:LEU:CB	1:A:368:MET:HE1	2.42	0.50
1:E:176:ARG:HH22	1:E:502:TYR:HB3	1.76	0.50
1:A:434:VAL:HG22	1:A:441:LYS:HG2	1.92	0.50
1:E:41:ARG:HD2	1:E:347:SER:HB2	1.94	0.50
1:E:383:GLY:O	1:E:387:LYS:HG3	2.11	0.50
1:A:228:LEU:HD12	1:A:230:MET:HE3	1.94	0.50
1:A:440:GLN:HE21	1:A:440:GLN:CA	2.25	0.50
1:E:221:ARG:O	1:E:225:ARG:HG3	2.12	0.50
1:E:179:GLU:OE1	1:E:501:GLN:HG3	2.11	0.50
1:C:81:ASP:OD2	1:C:82:PRO:CD	2.60	0.49
1:C:435:PRO:O	1:C:437:ASP:N	2.45	0.49
1:C:118:ILE:HG22	1:C:119:ARG:N	2.28	0.49
1:E:51:GLU:O	1:E:51:GLU:HG2	2.12	0.49
1:A:261:THR:O	1:A:327:SER:OG	2.30	0.49
1:C:457:LYS:HE3	1:C:460:ARG:HH12	1.78	0.49
1:E:123:GLY:N	1:E:128:ARG:O	2.43	0.49
1:E:86:LEU:HD22	1:E:293:TRP:HA	1.95	0.49
1:E:219:LEU:O	1:E:223:MET:HG3	2.13	0.49
1:E:122:LEU:HA	1:E:128:ARG:O	2.13	0.49
1:E:226:GLN:H	1:E:227:GLY:HA2	1.77	0.49
1:C:39:GLU:CG	1:C:41:ARG:HH21	2.15	0.49
1:C:86:LEU:HD22	1:C:293:TRP:HA	1.95	0.48
1:C:219:LEU:O	1:C:223:MET:HG3	2.12	0.48
1:C:109:ASP:HA	1:C:183:THR:HB	1.95	0.48
1:A:163:VAL:HG21	1:A:230:MET:HE1	1.96	0.48
1:E:437:ASP:OD2	1:E:438:GLY:HA3	2.13	0.48
1:C:179:GLU:HA	1:C:231:LYS:O	2.13	0.48
1:C:179:GLU:OE1	1:C:501:GLN:HB3	2.13	0.48
1:C:263:THR:HG22	1:C:323:GLU:HG3	1.95	0.48
1:E:433:ARG:C	1:E:434:VAL:CG1	2.87	0.48
1:C:378:ASP:OD1	1:C:378:ASP:N	2.47	0.48
1:A:118:ILE:HD13	1:A:489:GLY:HA3	1.95	0.48
1:A:219:LEU:O	1:A:223:MET:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:LYS:HE3	1:A:380:LYS:HB3	1.60	0.48
1:C:50:GLU:O	1:C:54:GLN:HG2	2.14	0.48
1:C:435:PRO:C	1:C:437:ASP:N	2.72	0.48
1:E:218:GLU:OE2	1:E:221:ARG:NH2	2.47	0.47
1:C:437:ASP:CG	1:C:438:GLY:HA3	2.36	0.47
1:A:264:ASN:HA	1:A:293:TRP:CD1	2.50	0.47
1:E:70:MET:O	1:E:74:MET:HG3	2.14	0.47
1:C:106:TYR:CZ	1:C:119:ARG:HD3	2.49	0.47
1:E:431:ARG:HD2	1:E:444:THR:HG21	1.96	0.47
1:C:225:ARG:O	1:E:83:HIS:NE2	2.35	0.47
1:C:341:LYS:HE3	1:C:345:ASP:OD2	2.15	0.47
1:C:264:ASN:HA	1:C:293:TRP:CD1	2.49	0.47
1:E:332:GLY:N	1:E:365:THR:HG22	2.27	0.47
1:C:70:MET:O	1:C:74:MET:HG3	2.14	0.47
1:E:435:PRO:CA	1:E:436:SER:CB	2.93	0.47
1:C:80:ALA:HB2	1:C:276:LYS:CB	2.44	0.47
1:A:303:ARG:NH2	1:A:308:ASN:OD1	2.48	0.47
1:A:140:PRO:HD2	1:A:143:LEU:HD12	1.97	0.47
1:C:434:VAL:HG21	1:C:474:GLU:CD	2.40	0.46
1:A:59:PRO:O	1:A:62:LEU:N	2.47	0.46
1:E:474:GLU:O	1:E:477:SER:OG	2.32	0.46
1:E:105:PHE:CZ	1:E:497:ALA:HA	2.51	0.46
1:A:180:LEU:HD23	1:A:232:VAL:CG2	2.42	0.46
1:C:433:ARG:HA	1:C:433:ARG:HD3	1.56	0.46
1:C:160:GLU:OE2	1:C:164:LYS:HE3	2.16	0.46
1:E:88:MET:HE3	1:E:289:ILE:HG21	1.98	0.46
1:E:173:PRO:HG2	1:E:176:ARG:HH11	1.81	0.46
1:C:247:ARG:HA	1:C:250:ASP:O	2.15	0.46
1:E:141:PRO:HA	1:E:144:MET:CE	2.46	0.46
1:A:91:SER:H	1:A:287:MET:HE2	1.81	0.46
1:C:380:LYS:HB2	1:C:380:LYS:HE3	1.61	0.46
1:E:153:ASP:OD1	1:E:225:ARG:NH1	2.49	0.46
1:A:42:ARG:HE	1:A:42:ARG:HB2	1.56	0.46
1:A:391:GLY:O	1:C:352:VAL:HG11	2.16	0.45
1:C:387:LYS:HA	1:C:392:VAL:HG22	1.98	0.45
1:C:332:GLY:H	1:C:365:THR:CG2	2.24	0.45
1:E:65:GLY:O	1:E:300:ARG:NH1	2.49	0.45
1:C:173:PRO:HG2	1:C:176:ARG:NH1	2.30	0.45
1:A:109:ASP:O	1:A:115:PHE:HA	2.16	0.45
1:E:298:SER:C	1:E:300:ARG:H	2.25	0.45
1:C:41:ARG:C	1:C:43:ARG:N	2.75	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:PHE:CZ	1:C:497:ALA:HA	2.52	0.45
1:C:329:MET:HE2	1:C:330:TYR:CE1	2.51	0.45
1:E:438:GLY:N	1:E:441:LYS:HZ3	2.14	0.45
1:C:244:ALA:O	1:C:245:GLY:C	2.57	0.45
1:E:435:PRO:HB3	1:E:436:SER:HB2	1.99	0.45
1:E:440:GLN:CA	1:E:441:LYS:CB	2.39	0.45
1:A:472:GLY:C	1:A:474:GLU:N	2.74	0.45
1:E:120:VAL:HG12	1:E:132:GLN:HG3	1.98	0.45
1:E:333:GLU:O	1:E:337:ARG:HG3	2.17	0.45
1:A:183:THR:HA	1:A:236:VAL:O	2.17	0.44
1:E:39:GLU:OE2	1:E:40:GLU:N	2.47	0.44
1:C:173:PRO:CG	1:C:176:ARG:HH11	2.27	0.44
1:E:296:PHE:O	1:E:322:TYR:HB2	2.17	0.44
1:E:95:ASN:HD21	1:E:283:ARG:CD	2.30	0.44
1:E:109:ASP:HA	1:E:183:THR:HB	1.99	0.44
1:A:483:LEU:CD1	1:A:485:ASN:HB2	2.48	0.44
1:E:431:ARG:HB3	1:E:444:THR:HG21	2.00	0.44
1:C:80:ALA:HB2	1:C:276:LYS:HB2	1.99	0.44
1:A:226:GLN:N	1:A:227:GLY:HA2	2.33	0.44
1:C:277:TRP:CZ3	1:C:287:MET:HE2	2.52	0.44
1:E:155:ILE:HD13	1:E:155:ILE:HG21	1.78	0.44
1:C:216:VAL:CG1	1:C:232:VAL:HG11	2.46	0.43
1:C:434:VAL:HG21	1:C:474:GLU:OE1	2.16	0.43
1:E:216:VAL:HG23	1:E:235:LEU:HD22	1.98	0.43
1:E:287:MET:HE3	1:E:288:VAL:N	2.28	0.43
1:E:440:GLN:HA	1:E:441:LYS:HB2	0.60	0.43
1:A:166:GLU:CD	1:A:178:ARG:HH12	2.26	0.43
1:E:264:ASN:HB2	2:E:601:BGC:H5	1.99	0.43
1:A:160:GLU:HB2	1:A:228:LEU:HD22	2.00	0.43
1:A:447:ALA:C	1:A:448:LEU:HD12	2.43	0.43
1:E:298:SER:C	1:E:300:ARG:N	2.76	0.43
1:E:437:ASP:HA	1:E:438:GLY:HA3	1.79	0.43
1:A:500:SER:C	1:A:502:TYR:H	2.26	0.43
1:C:168:GLU:OE2	1:C:168:GLU:HA	2.18	0.43
1:A:296:PHE:HD2	1:A:322:TYR:CD2	2.37	0.43
1:E:335:VAL:HG12	1:E:410:VAL:HG21	1.99	0.43
1:A:296:PHE:O	1:A:322:TYR:HB2	2.18	0.43
1:C:118:ILE:CG2	1:C:119:ARG:N	2.82	0.43
1:C:328:GLY:O	1:C:365:THR:HG21	2.18	0.43
1:E:368:MET:HE3	1:E:456:TYR:CE2	2.45	0.43
1:C:88:MET:HE3	1:C:88:MET:HB3	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:GLY:O	1:C:207:ILE:HG23	2.19	0.43
1:C:247:ARG:NH2	1:C:253:VAL:O	2.46	0.43
1:C:427:LYS:HE2	1:C:427:LYS:HB3	1.83	0.43
1:E:264:ASN:CG	1:E:265:ALA:H	2.27	0.43
1:E:436:SER:N	1:E:437:ASP:CA	2.81	0.43
1:C:44:ARG:HD3	1:C:345:ASP:O	2.18	0.43
1:C:448:LEU:HD13	1:C:481:VAL:HG13	2.01	0.43
1:E:264:ASN:CG	1:E:265:ALA:N	2.76	0.43
1:A:396:SER:O	1:A:400:ARG:HG3	2.19	0.42
1:C:447:ALA:C	1:C:448:LEU:HD12	2.44	0.42
1:A:68:ASP:OD1	1:A:427:LYS:NZ	2.46	0.42
1:E:323:GLU:OE2	2:E:601:BGC:O1	2.37	0.42
1:C:293:TRP:CD1	1:C:293:TRP:C	2.97	0.42
1:A:252:ASP:HB3	1:A:445:VAL:HG23	2.01	0.42
1:C:389:ILE:HG22	1:C:390:LEU:HD23	2.01	0.42
1:A:106:TYR:CB	1:A:180:LEU:HD12	2.49	0.42
1:C:220:SER:O	1:C:224:GLU:HG3	2.20	0.42
1:E:184:PHE:CE2	1:E:186:PHE:HB2	2.54	0.42
1:A:397:LEU:HD13	1:A:400:ARG:HH11	1.85	0.42
1:E:128:ARG:NH1	1:E:248:TYR:O	2.53	0.42
1:C:45:ALA:HB2	1:C:346:ALA:CB	2.50	0.42
1:E:58:THR:HB	1:E:63:LEU:HD21	2.02	0.42
1:E:118:ILE:HG22	1:E:119:ARG:N	2.34	0.42
1:E:435:PRO:CA	1:E:436:SER:HB2	2.50	0.42
1:C:497:ALA:C	1:C:499:HIS:H	2.28	0.42
1:E:261:THR:C	1:E:327:SER:HG	2.26	0.42
1:C:43:ARG:O	1:C:46:ALA:HB3	2.20	0.42
1:E:373:HIS:CD2	1:E:458:LYS:HG3	2.54	0.41
1:A:225:ARG:O	1:C:83:HIS:NE2	2.51	0.41
1:E:293:TRP:CD1	1:E:293:TRP:C	2.98	0.41
1:C:457:LYS:CD	1:C:460:ARG:HH12	2.32	0.41
1:A:176:ARG:HD3	1:A:503:ALA:CA	2.49	0.41
1:A:435:PRO:HB3	1:A:436:SER:HB2	2.02	0.41
1:E:257:VAL:HB	1:E:448:LEU:HD13	2.03	0.41
1:E:438:GLY:N	1:E:441:LYS:HZ1	2.14	0.41
1:C:264:ASN:CG	1:C:265:ALA:H	2.28	0.41
1:A:67:ALA:O	1:A:71:VAL:HG23	2.20	0.41
1:A:102:HIS:HA	1:A:122:LEU:HB2	2.03	0.41
1:A:437:ASP:OD2	1:A:439:SER:CB	2.68	0.41
1:C:126:GLU:HG2	1:C:127:LYS:N	2.35	0.41
1:E:243:LEU:HB2	1:E:256:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:SER:HA	1:A:164:LYS:HG2	2.03	0.41
1:A:201:TRP:HE1	1:A:210:THR:HG23	1.85	0.41
1:E:73:GLU:OE2	1:E:76:ARG:HD3	2.20	0.41
1:E:501:GLN:N	1:E:502:TYR:HA	2.35	0.41
1:A:83:HIS:NE2	1:E:225:ARG:O	2.40	0.41
1:A:437:ASP:OD1	1:A:437:ASP:N	2.54	0.41
1:A:475:ALA:O	1:A:478:SER:OG	2.38	0.41
1:A:497:ALA:C	1:A:499:HIS:N	2.79	0.41
1:C:386:LEU:O	1:C:390:LEU:HB2	2.20	0.41
1:C:445:VAL:HG22	1:C:480:VAL:HB	2.03	0.41
1:E:94:ASP:OD1	1:E:94:ASP:N	2.45	0.41
1:E:176:ARG:NH2	1:E:502:TYR:C	2.62	0.41
1:C:39:GLU:CD	1:C:41:ARG:NH2	2.75	0.41
1:E:389:ILE:HG22	1:E:390:LEU:HD23	2.02	0.41
1:E:378:ASP:OD1	1:E:378:ASP:N	2.47	0.40
1:A:49:ILE:O	1:A:53:GLU:CG	2.68	0.40
1:A:60:THR:O	1:A:64:ARG:HG3	2.22	0.40
1:A:176:ARG:HD3	1:A:503:ALA:CB	2.51	0.40
1:A:277:TRP:NE1	1:A:281:LEU:HD23	2.37	0.40
1:A:485:ASN:C	1:A:487:GLY:N	2.77	0.40
1:E:247:ARG:NH2	1:E:253:VAL:O	2.47	0.40
1:A:164:LYS:HB3	1:A:164:LYS:HE3	1.90	0.40
1:C:397:LEU:HA	1:C:400:ARG:HG3	2.02	0.40
1:E:108:LEU:HD21	1:E:155:ILE:HG21	2.03	0.40
1:E:143:LEU:HA	1:E:143:LEU:HD23	1.92	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	463/465 (100%)	435 (94%)	25 (5%)	3 (1%)	21 51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	463/465 (100%)	435 (94%)	22 (5%)	6 (1%)	9	32
1	E	463/465 (100%)	438 (95%)	23 (5%)	2 (0%)	30	59
All	All	1389/1395 (100%)	1308 (94%)	70 (5%)	11 (1%)	16	44

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	436	SER
1	E	441	LYS
1	A	436	SER
1	C	42	ARG
1	C	436	SER
1	C	440	GLN
1	A	42	ARG
1	A	439	SER
1	C	41	ARG
1	C	500	SER
1	C	435	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	377/377 (100%)	344 (91%)	33 (9%)	9	29
1	C	377/377 (100%)	355 (94%)	22 (6%)	18	49
1	E	377/377 (100%)	362 (96%)	15 (4%)	28	62
All	All	1131/1131 (100%)	1061 (94%)	70 (6%)	16	46

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ARG
1	A	53	GLU
1	A	55	ARG

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Mol	Chain	Res	Type
1	A	75	GLU
1	A	116	ARG
1	A	117	VAL
1	A	126	GLU
1	A	131	SER
1	A	164	LYS
1	A	168	GLU
1	A	172	LEU
1	A	173	PRO
1	A	197	THR
1	A	251	ASN
1	A	252	ASP
1	A	261	THR
1	A	293	TRP
1	A	351	ASP
1	A	363	LEU
1	A	365	THR
1	A	375	THR
1	A	380	LYS
1	A	437	ASP
1	A	439	SER
1	A	440	GLN
1	A	441	LYS
1	A	444	THR
1	A	473	GLU
1	A	474	GLU
1	A	483	LEU
1	A	488	SER
1	A	490	ILE
1	A	501	GLN
1	C	41	ARG
1	C	44	ARG
1	C	50	GLU
1	C	116	ARG
1	C	117	VAL
1	C	132	GLN
1	C	197	THR
1	C	243	LEU
1	C	287	MET
1	C	293	TRP
1	C	297	LYS
1	C	363	LEU

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Mol	Chain	Res	Type
1	C	387	LYS
1	C	388	ASP
1	C	394	ASP
1	C	434	VAL
1	C	436	SER
1	C	437	ASP
1	C	439	SER
1	C	440	GLN
1	C	442	GLN
1	C	501	GLN
1	E	104	LEU
1	E	128	ARG
1	E	131	SER
1	E	173	PRO
1	E	176	ARG
1	E	283	ARG
1	E	293	TRP
1	E	335	VAL
1	E	365	THR
1	E	380	LYS
1	E	398	GLU
1	E	436	SER
1	E	441	LYS
1	E	442	GLN
1	E	444	THR

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	226	GLN
1	A	377	HIS
1	A	440	GLN
1	C	132	GLN
1	C	189	HIS
1	C	373	HIS
1	E	142	HIS
1	E	189	HIS
1	E	286	ASN
1	E	320	GLN
1	E	359	GLN
1	E	442	GLN
1	E	501	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	BGC	C	601	-	12,12,12	1.15	2 (16%)	17,17,17	1.98	3 (17%)
2	BGC	E	601	-	12,12,12	1.22	1 (8%)	17,17,17	0.83	0
2	BGC	A	601	-	12,12,12	1.40	2 (16%)	17,17,17	1.59	3 (17%)

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	BGC	C	601	-	-	2/2/22/22	0/1/1/1
2	BGC	E	601	-	-	0/2/22/22	0/1/1/1
2	BGC	A	601	-	-	2/2/22/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	601	BGC	O5-C1	3.20	1.50	1.42
2	A	601	BGC	C3-C2	-2.53	1.45	1.52
2	A	601	BGC	O1-C1	-2.27	1.32	1.39
2	C	601	BGC	O5-C1	2.14	1.48	1.42
2	C	601	BGC	O1-C1	-2.04	1.33	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	BGC	C1-C2-C3	-5.08	100.00	110.36
2	C	601	BGC	C1-O5-C5	-4.45	105.04	113.65
2	A	601	BGC	C1-C2-C3	-3.57	103.06	110.36
2	A	601	BGC	C1-O5-C5	-2.61	108.60	113.65
2	A	601	BGC	O4-C4-C3	-2.50	104.48	110.38
2	C	601	BGC	O5-C1-C2	-2.13	106.56	110.30

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	BGC	O5-C5-C6-O6
2	C	601	BGC	C4-C5-C6-O6
2	A	601	BGC	C4-C5-C6-O6
2	A	601	BGC	O5-C5-C6-O6

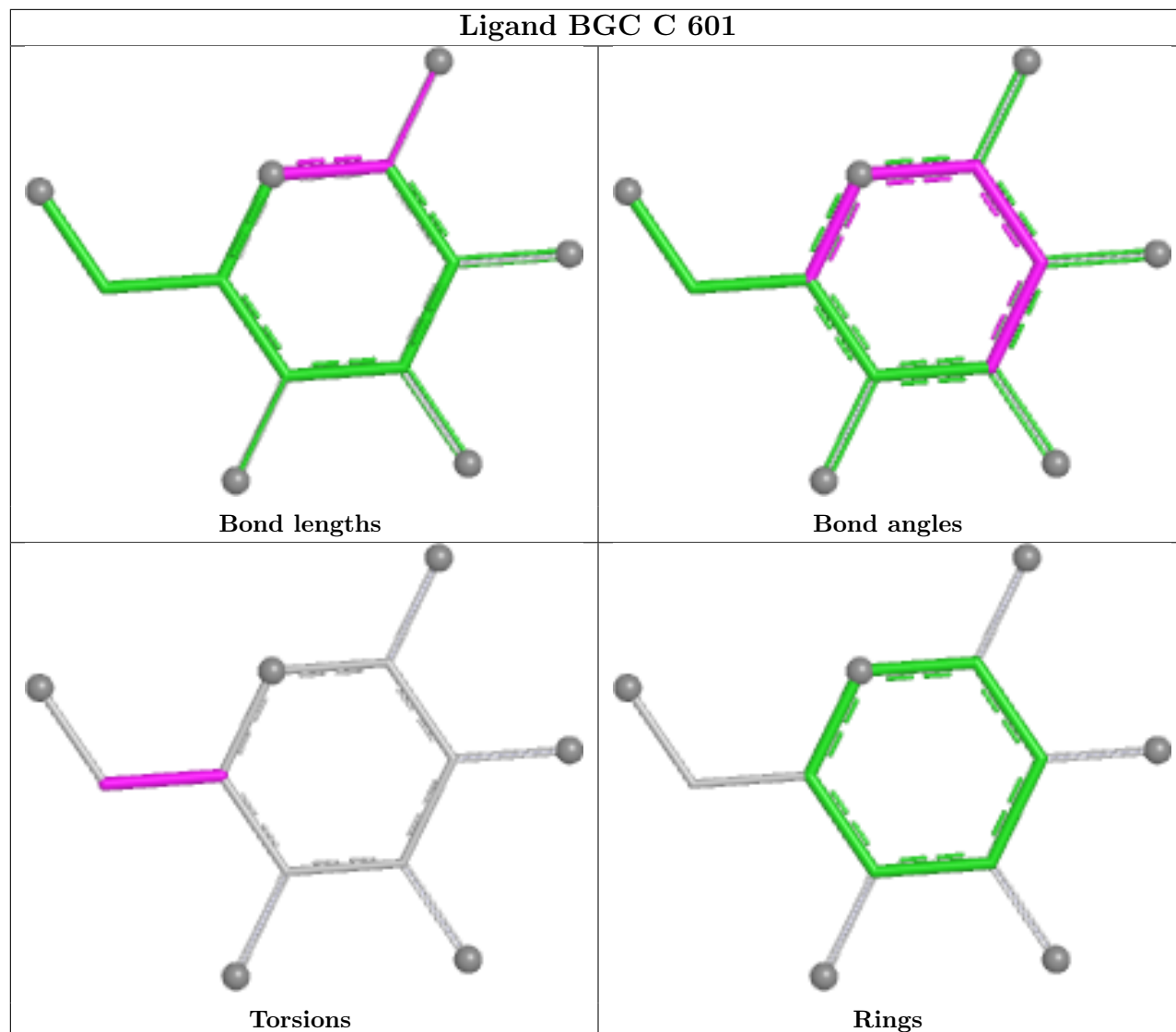
There are no ring outliers.

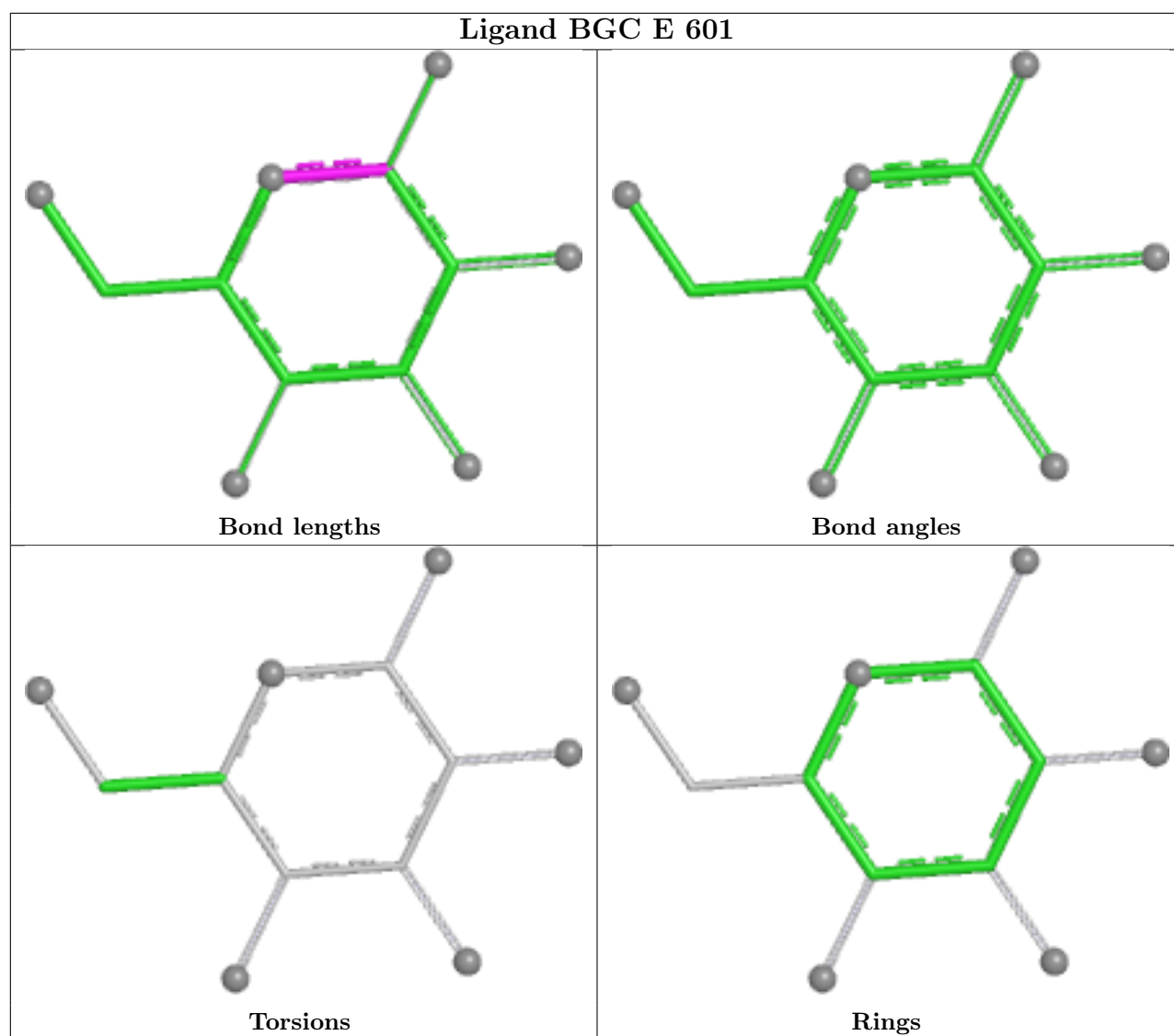
1 monomer is involved in 2 short contacts:

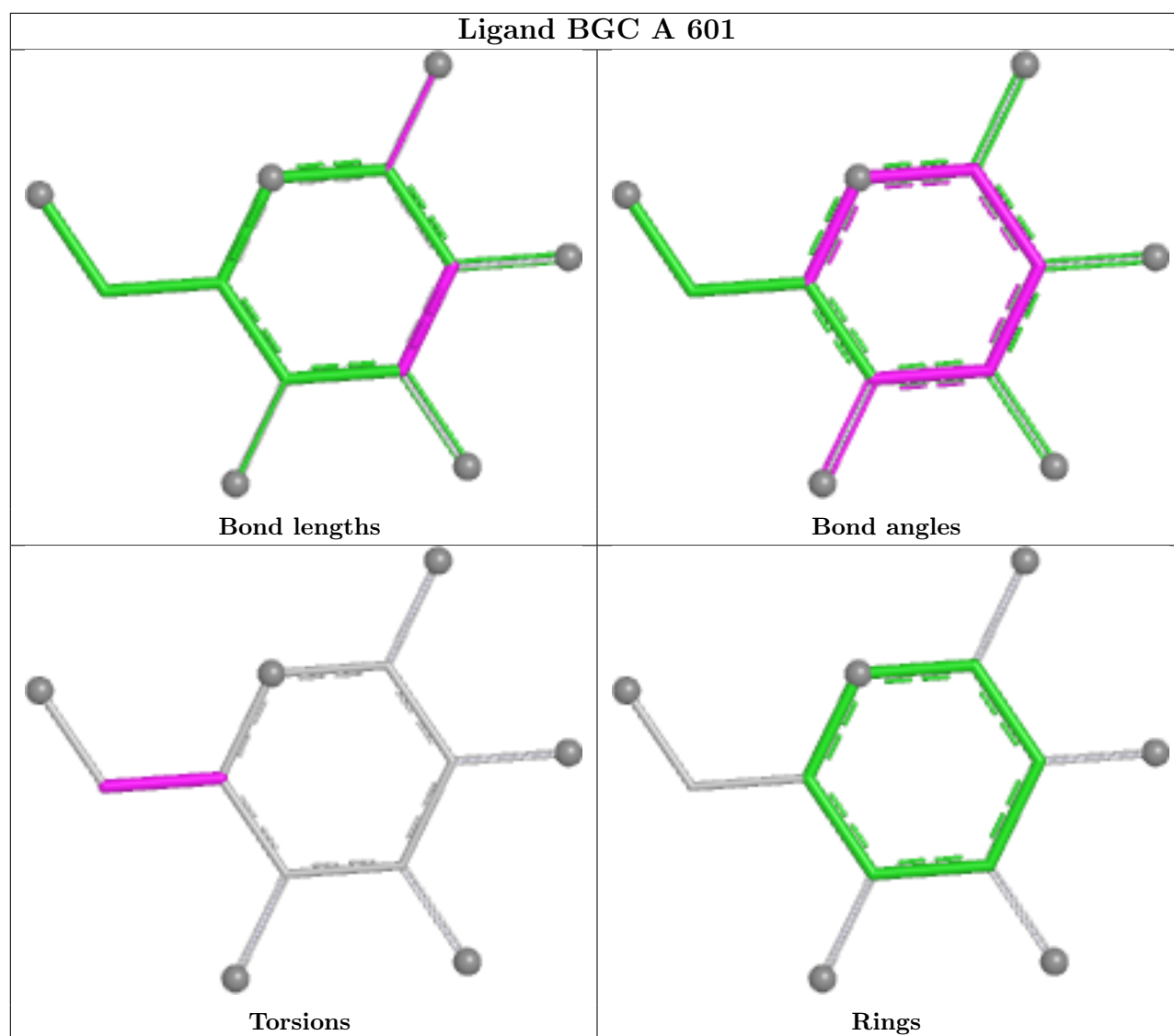
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	601	BGC	2	0

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	465/465 (100%)	-0.23	5 (1%) 78 71	14, 24, 49, 89	0
1	C	465/465 (100%)	0.07	11 (2%) 59 50	14, 34, 67, 122	0
1	E	465/465 (100%)	-0.13	11 (2%) 59 50	13, 25, 53, 118	0
All	All	1395/1395 (100%)	-0.09	27 (1%) 66 58	13, 27, 60, 122	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	435	PRO	5.5
1	C	485	ASN	3.9
1	C	441	LYS	3.7
1	C	502	TYR	3.6
1	E	94	ASP	3.4
1	E	434	VAL	3.4
1	E	436	SER	3.3
1	A	502	TYR	3.2
1	E	502	TYR	3.2
1	E	438	GLY	3.2
1	E	441	LYS	2.9
1	E	439	SER	2.9
1	C	377	HIS	2.8
1	C	175	GLY	2.8
1	E	503	ALA	2.7
1	C	474	GLU	2.6
1	C	440	GLN	2.5
1	C	486	ASP	2.5
1	E	437	ASP	2.4
1	C	439	SER	2.4
1	A	437	ASP	2.3
1	A	436	SER	2.3
1	A	174	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	377	HIS	2.2
1	C	434	VAL	2.2
1	E	486	ASP	2.1
1	C	435	PRO	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](#)

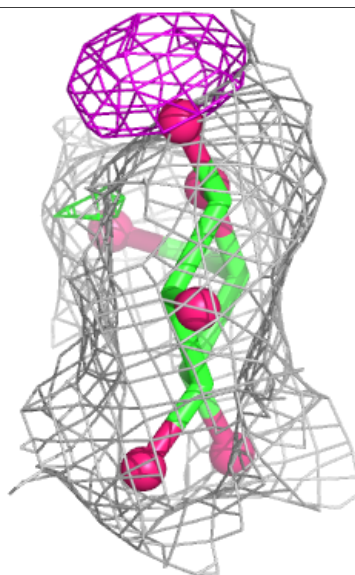
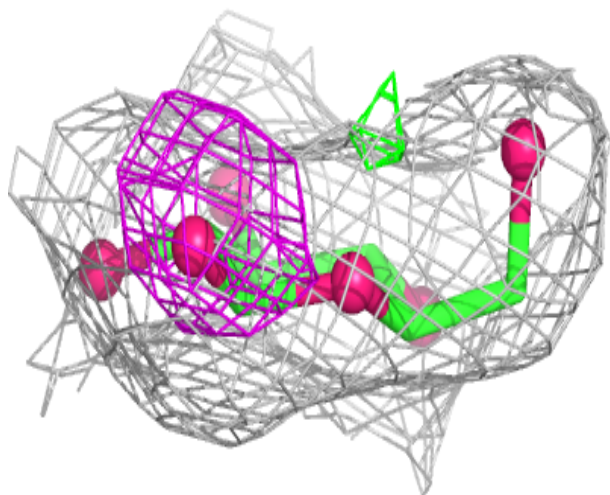
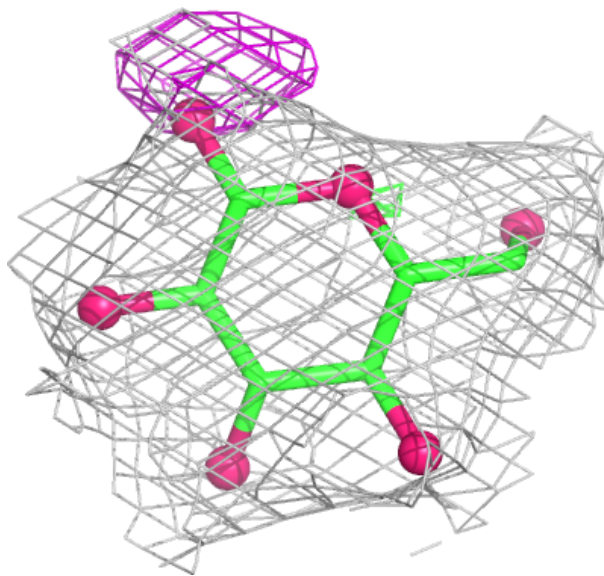
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	BGC	E	601	12/12	0.89	0.13	15,16,17,17	0
2	BGC	C	601	12/12	0.95	0.08	20,22,23,23	0
2	BGC	A	601	12/12	0.98	0.06	15,16,17,17	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.

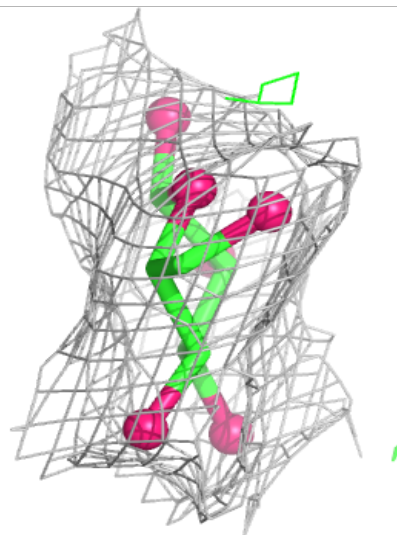
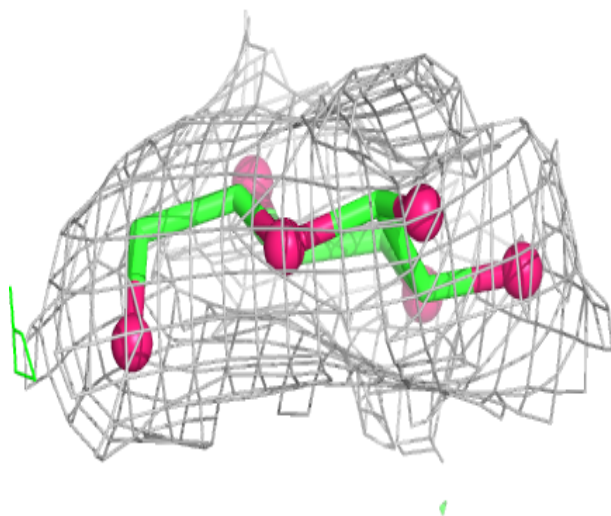
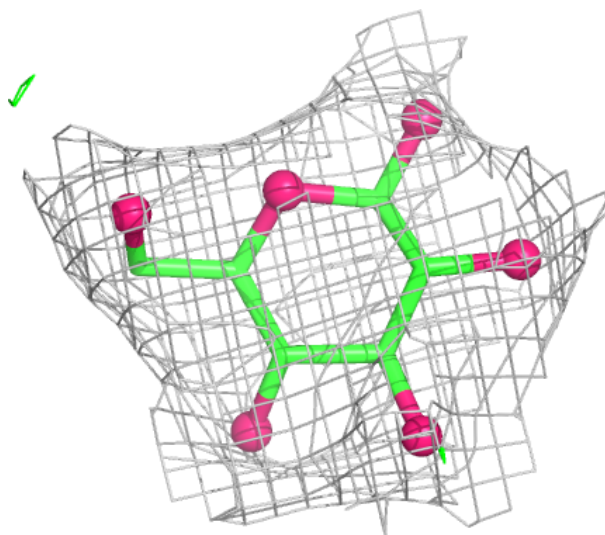
Electron density around BGC E 601:

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



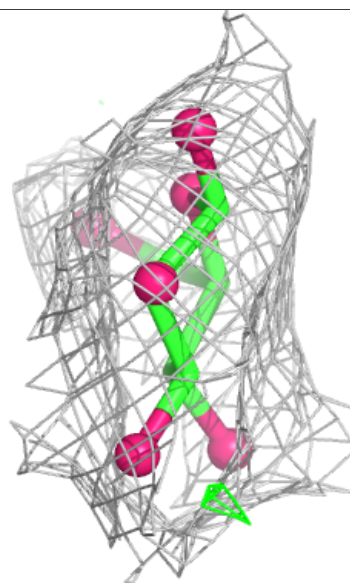
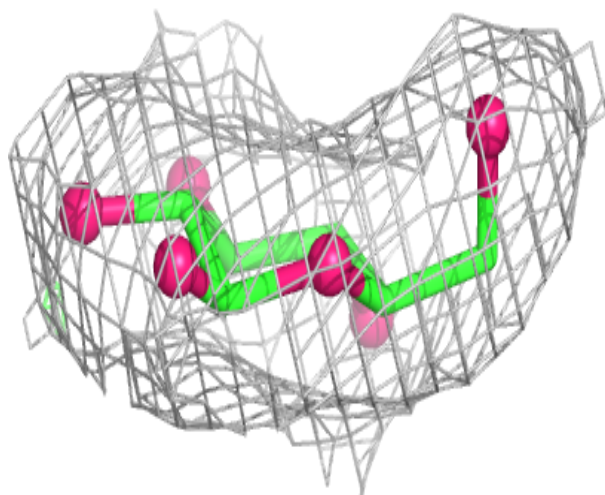
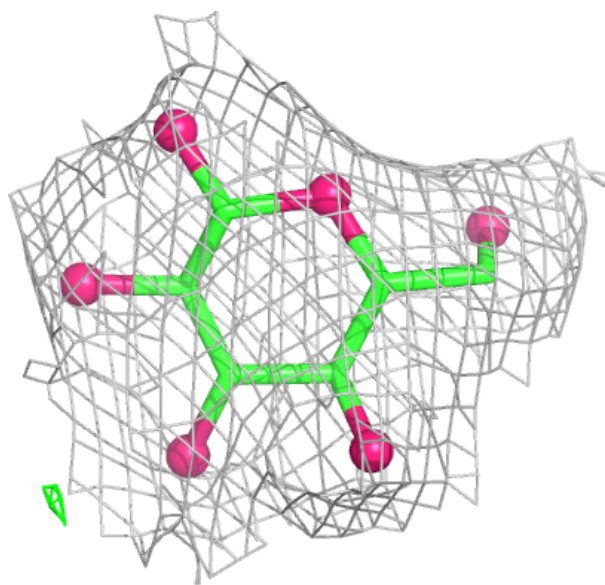
Electron density around BGC C 601:

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



Electron density around BGC A 601:

$2mF_o-DF_c$ (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.