



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 08:16 PM UTC

PDB ID : 8XRV / pdb_00008xrv
Title : The crystal structure of a GH3 enzyme CcBgl3B with glucose
Authors : Su, J.Y.
Deposited on : 2024-01-08
Resolution : 2.40 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:

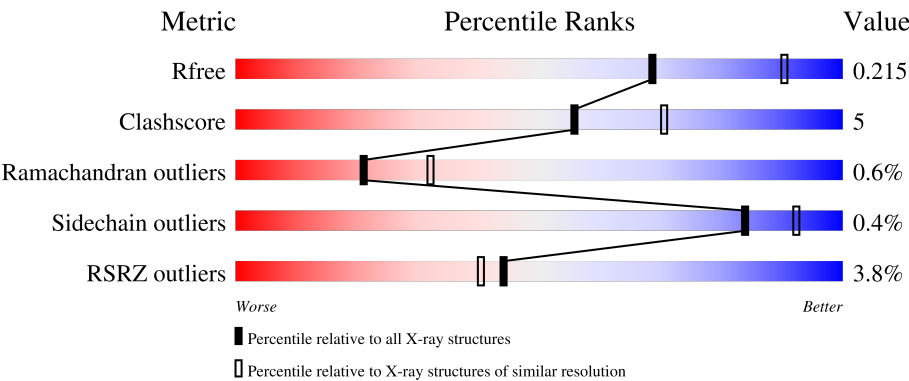
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	768	0% 87% 10% .
1	B	768	2% 84% 13% .
1	C	768	3% 88% 10% .
1	D	768	2% 89% 8% .
1	E	768	11% 82% 14% .

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Mol	Chain	Length	Quality of chain
1	F	768	% 88% 9%

2 Entry composition [i](#)

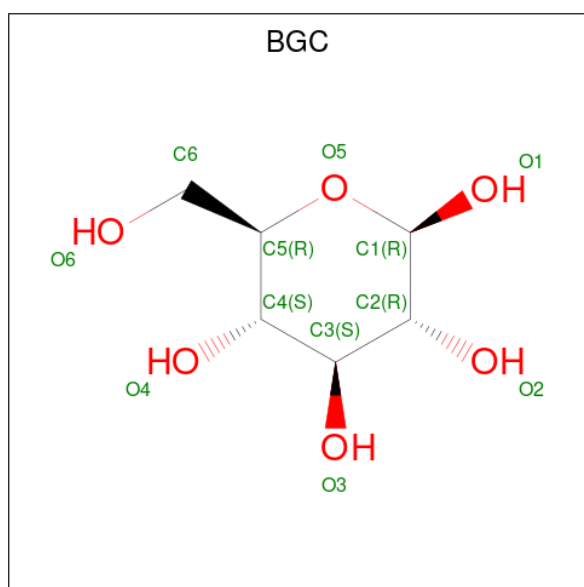
There are 4 unique types of molecules in this entry. The entry contains 35873 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GH3 enzyme CcBgl3B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	749	Total	C	N	O	S	0	0	0
			5623	3541	997	1070	15			
1	B	746	Total	C	N	O	S	0	0	0
			5596	3524	994	1063	15			
1	C	747	Total	C	N	O	S	0	0	0
			5609	3530	996	1068	15			
1	F	747	Total	C	N	O	S	0	0	0
			5598	3524	995	1064	15			
1	D	748	Total	C	N	O	S	0	0	0
			5605	3530	996	1064	15			
1	E	740	Total	C	N	O	S	0	0	0
			5547	3491	986	1055	15			

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0
3	B	1	Total Ca 1 1	0	0
3	C	1	Total Ca 1 1	0	0
3	F	1	Total Ca 1 1	0	0
3	D	1	Total Ca 1 1	0	0
3	E	1	Total Ca 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	391	Total O 391 391	0	0
4	B	360	Total O 360 360	0	0
4	C	398	Total O 398 398	0	0
4	F	447	Total O 447 447	0	0
4	D	353	Total O 353 353	0	0

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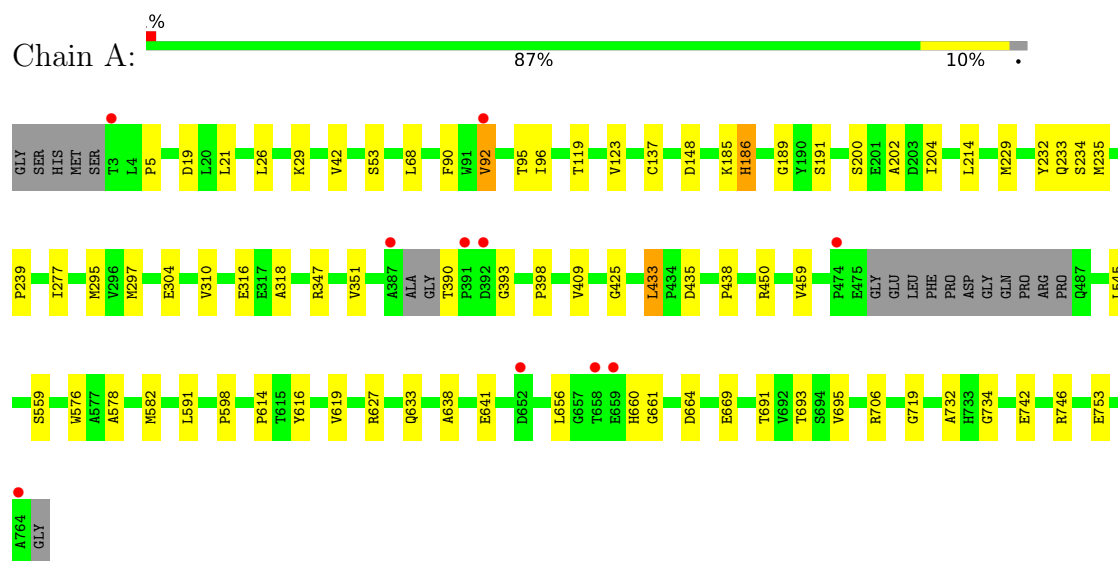
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	268	Total 268	O 268	0	0

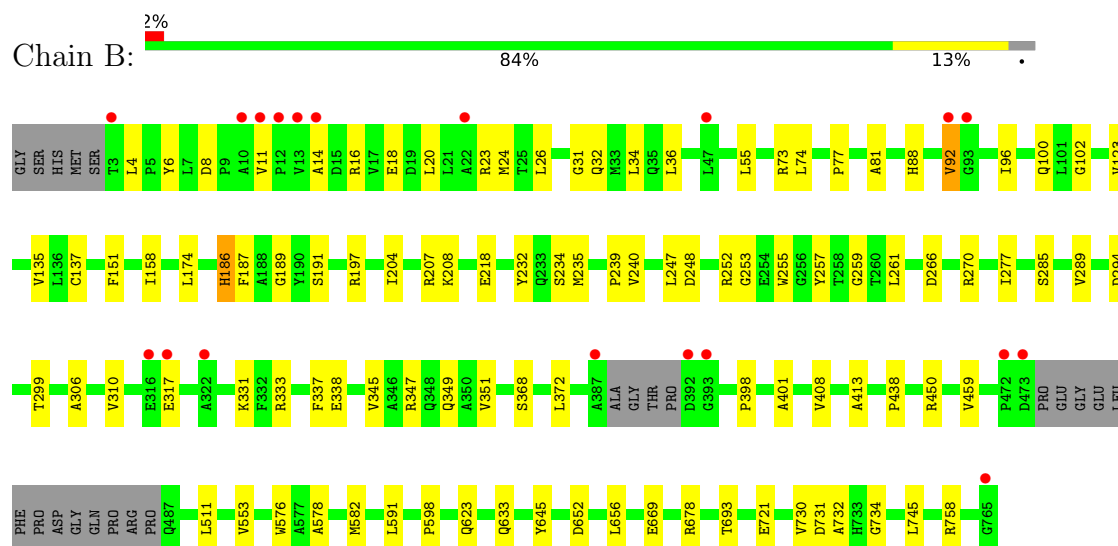
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

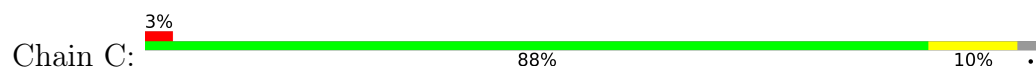
• Molecule 1: GH3 enzyme CcBgl3B

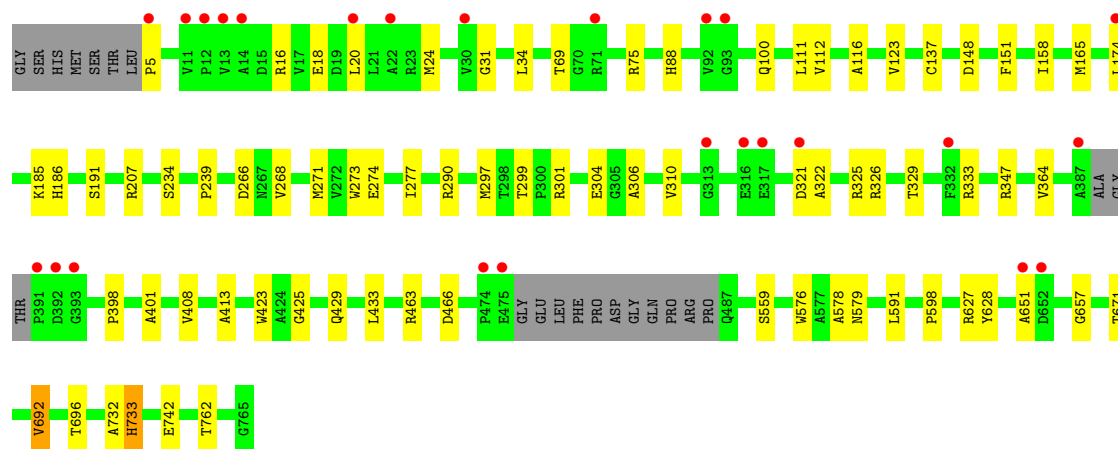


• Molecule 1: GH3 enzyme CcBgl3B

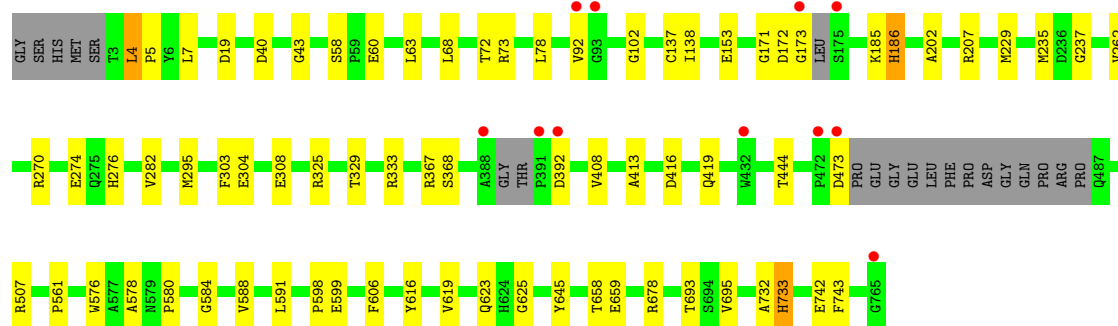
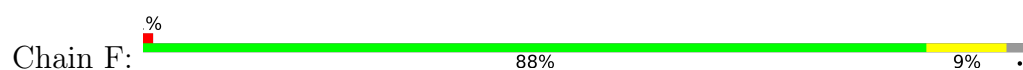


• Molecule 1: GH3 enzyme CcBgl3B

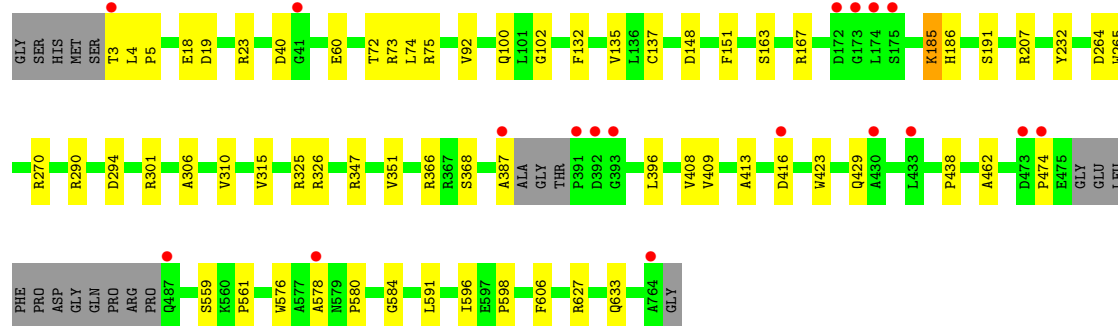
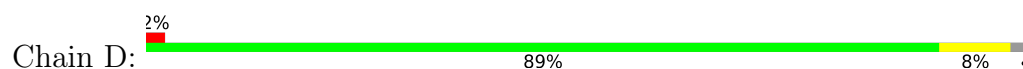




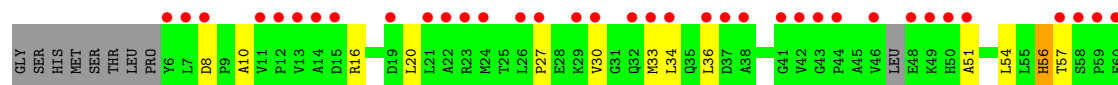
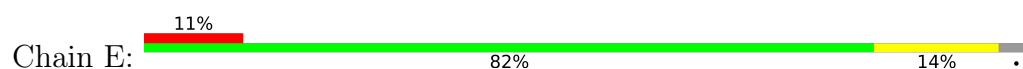
• Molecule 1: GH3 enzyme CcBgl3B

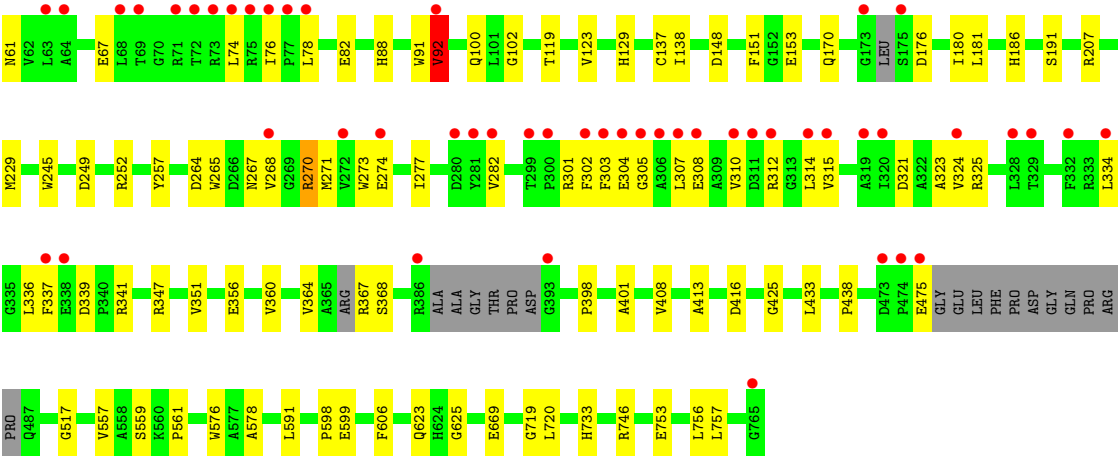


• Molecule 1: GH3 enzyme CcBgl3B



• Molecule 1: GH3 enzyme CcBgl3B





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.33Å 177.57Å 265.79Å 90.00° 100.86° 90.00°	Depositor
Resolution (Å)	19.97 – 2.40 19.97 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.97-2.40) 99.4 (19.97-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.41Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.176 , 0.215 0.178 , 0.215	Depositor DCC
R_{free} test set	2000 reflections (0.69%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.432	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	35873	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% 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: BGC, CA

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	0.40	0/5746	0.65	3/7859 (0.0%)
1	B	0.41	0/5717	0.66	1/7816 (0.0%)
1	C	0.42	0/5732	0.67	2/7837 (0.0%)
1	D	0.41	0/5727	0.66	2/7832 (0.0%)
1	E	0.38	0/5665	0.64	0/7741
1	F	0.42	0/5718	0.68	2/7816 (0.0%)
All	All	0.41	0/34305	0.66	10/46901 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	185	LYS	CA-C-N	5.82	132.65	121.54
1	C	185	LYS	C-N-CA	5.82	132.65	121.54
1	D	185	LYS	CA-C-N	5.38	131.82	121.54
1	D	185	LYS	C-N-CA	5.38	131.82	121.54
1	A	92	VAL	CA-C-O	-5.19	114.29	120.78
1	A	185	LYS	CA-C-N	5.16	131.39	121.54
1	A	185	LYS	C-N-CA	5.16	131.39	121.54
1	B	721	GLU	CA-CB-CG	5.11	124.33	114.10
1	F	185	LYS	CA-C-N	5.08	131.24	121.54
1	F	185	LYS	C-N-CA	5.08	131.24	121.54

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	5623	0	5536	53	0
1	B	5596	0	5512	68	0
1	C	5609	0	5519	54	0
1	D	5605	0	5508	40	0
1	E	5547	0	5445	65	0
1	F	5598	0	5507	42	0
2	A	12	0	12	2	0
2	B	12	0	11	1	0
2	C	12	0	11	2	0
2	D	12	0	12	1	0
2	E	12	0	12	2	0
2	F	12	0	11	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	391	0	0	4	0
4	B	360	0	0	2	0
4	C	398	0	0	3	0
4	D	353	0	0	3	0
4	E	268	0	0	2	0
4	F	447	0	0	6	0
All	All	35873	0	33096	316	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 5.

All (316) 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:96:ILE:H	1:A:582:MET:HE2	1.27	0.97
1:B:623:GLN:HE22	1:B:633:GLN:HE22	1.15	0.94
1:B:26:LEU:HD21	1:B:310:VAL:HG21	1.53	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:282:VAL:HG21	1:F:308:GLU:OE1	1.75	0.86
1:E:229:MET:HE1	2:E:801:BGC:H3	1.62	0.80
1:D:100:GLN:HG3	1:D:151:PHE:CE1	2.17	0.79
1:F:658:THR:HG23	1:F:659:GLU:HG2	1.69	0.73
1:A:691:THR:HB	1:A:742:GLU:HG3	1.71	0.72
1:C:329:THR:HG22	1:C:333:ARG:HE	1.52	0.72
1:E:271:MET:HG2	1:E:277:ILE:HD11	1.69	0.72
1:C:207:ARG:NH2	4:C:903:HOH:O	2.24	0.70
1:B:100:GLN:HG3	1:B:151:PHE:CE1	2.26	0.70
1:A:96:ILE:N	1:A:582:MET:HE2	2.06	0.70
1:E:207:ARG:NH2	4:E:901:HOH:O	2.23	0.70
1:A:53:SER:HB3	1:A:295:MET:HE1	1.74	0.70
1:E:100:GLN:HG3	1:E:151:PHE:CE1	2.26	0.70
1:D:301:ARG:HG2	1:D:301:ARG:HH11	1.57	0.69
1:B:591:LEU:HD11	1:B:598:PRO:HG3	1.75	0.68
1:F:5:PRO:HG2	1:F:19:ASP:CG	2.20	0.66
1:B:398:PRO:HG2	1:B:401:ALA:HB3	1.78	0.66
1:F:207:ARG:NH2	4:F:903:HOH:O	2.28	0.65
1:E:181:LEU:HD21	1:E:334:LEU:HD21	1.79	0.65
1:B:285:SER:O	1:B:289:VAL:HG23	1.97	0.65
1:B:266:ASP:HB3	1:B:270:ARG:HH21	1.64	0.63
1:C:347:ARG:NH1	4:C:909:HOH:O	2.31	0.63
1:D:265:TRP:O	1:D:270:ARG:NH2	2.32	0.62
1:B:137:CYS:HB3	1:B:191:SER:HB3	1.79	0.62
1:F:262:VAL:HG23	1:F:295:MET:HG2	1.82	0.61
1:B:207:ARG:NH1	1:B:730:VAL:O	2.33	0.61
1:C:325:ARG:HB2	1:C:326:ARG:HH11	1.66	0.61
1:E:367:ARG:NH1	1:E:599:GLU:OE1	2.33	0.61
1:E:252:ARG:NH2	1:E:257:TYR:O	2.27	0.60
1:B:34:LEU:HD13	1:B:36:LEU:HD11	1.83	0.60
1:D:207:ARG:NH2	4:D:902:HOH:O	2.21	0.60
1:B:189:GLY:HA3	1:B:235:MET:HE1	1.84	0.60
1:C:329:THR:HG21	1:C:333:ARG:HH21	1.65	0.60
1:B:232:TYR:OH	2:B:801:BGC:O1	2.20	0.60
1:C:591:LEU:HD11	1:C:598:PRO:HG3	1.82	0.60
1:A:576:TRP:CZ2	1:A:578:ALA:HB2	2.37	0.59
1:C:306:ALA:O	1:C:310:VAL:HG23	2.01	0.59
1:E:301:ARG:NH2	4:E:903:HOH:O	2.28	0.59
1:B:266:ASP:HB3	1:B:270:ARG:NH2	2.17	0.59
1:E:137:CYS:HB3	1:E:191:SER:HB3	1.84	0.59
1:B:174:LEU:HD13	1:B:333:ARG:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:322:ALA:O	1:C:326:ARG:NH1	2.36	0.59
1:F:591:LEU:HD11	1:F:598:PRO:HG3	1.85	0.58
1:C:277:ILE:HA	1:F:693:THR:O	2.04	0.58
1:C:651:ALA:HB3	1:C:671:THR:HB	1.86	0.58
1:B:32:GLN:OE1	1:B:331:LYS:NZ	2.38	0.57
1:C:88:HIS:CD2	1:C:123:VAL:HA	2.39	0.57
1:A:233:GLN:HE21	1:A:235:MET:HE3	1.68	0.57
1:B:88:HIS:CD2	1:B:123:VAL:HA	2.40	0.56
1:A:5:PRO:HG2	1:A:19:ASP:CG	2.30	0.56
1:A:627:ARG:HA	1:A:633:GLN:HG3	1.86	0.56
1:D:408:VAL:HG12	1:D:413:ALA:HB1	1.86	0.56
1:D:366:ARG:HG2	1:D:596:ILE:HD13	1.87	0.56
1:E:425:GLY:HA2	1:E:433:LEU:HD22	1.87	0.56
1:D:137:CYS:SG	1:D:186:HIS:HB2	2.46	0.56
1:F:202:ALA:HB3	1:F:235:MET:HG2	1.88	0.55
1:A:693:THR:O	1:B:277:ILE:HA	2.06	0.55
1:C:408:VAL:HG12	1:C:413:ALA:HB1	1.88	0.55
1:B:34:LEU:HB3	1:B:36:LEU:CD1	2.37	0.55
1:C:116:ALA:HB2	1:C:165:MET:HG2	1.88	0.55
1:C:273:TRP:CE3	1:C:274:GLU:HB2	2.42	0.55
1:D:100:GLN:HG3	1:D:151:PHE:CD1	2.42	0.55
1:F:60:GLU:CD	1:F:60:GLU:H	2.14	0.55
1:C:329:THR:O	1:C:333:ARG:HG3	2.06	0.55
1:F:137:CYS:SG	1:F:186:HIS:HB2	2.47	0.54
1:F:137:CYS:SG	4:F:1236:HOH:O	2.27	0.54
1:E:267:ASN:O	1:E:270:ARG:HB2	2.08	0.54
1:E:273:TRP:CZ3	1:E:274:GLU:HB2	2.42	0.54
1:D:294:ASP:OD1	1:D:326:ARG:NH1	2.40	0.54
1:A:137:CYS:SG	4:A:1184:HOH:O	2.26	0.54
1:B:338:GLU:N	1:B:338:GLU:OE2	2.40	0.54
1:E:229:MET:HE3	1:E:264:ASP:CG	2.33	0.54
1:B:576:TRP:CZ2	1:B:578:ALA:HB2	2.43	0.54
1:B:678:ARG:NH2	4:B:902:HOH:O	2.31	0.54
1:C:463:ARG:NH2	1:C:466:ASP:OD1	2.41	0.53
1:A:591:LEU:HD11	1:A:598:PRO:HG3	1.89	0.53
1:C:20:LEU:O	1:C:24:MET:HG3	2.08	0.53
1:F:561:PRO:HB3	1:F:606:PHE:CD1	2.43	0.53
1:B:100:GLN:NE2	1:B:135:VAL:O	2.41	0.53
1:A:706:ARG:NH1	1:A:719:GLY:O	2.41	0.53
1:C:297:MET:HE1	2:C:801:BGC:H6C2	1.89	0.53
1:D:100:GLN:NE2	1:D:135:VAL:O	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:282:VAL:HG22	1:E:305:GLY:HA2	1.88	0.53
1:A:42:VAL:HG12	1:A:68:LEU:HD12	1.90	0.53
1:A:390:THR:HG22	1:A:393:GLY:H	1.72	0.53
1:D:75:ARG:NH2	4:D:904:HOH:O	2.24	0.53
1:C:692:VAL:HG22	1:F:276:HIS:HB3	1.91	0.53
1:C:576:TRP:CZ2	1:C:578:ALA:HB2	2.43	0.53
1:D:591:LEU:HD11	1:D:598:PRO:HG3	1.91	0.53
1:A:656:LEU:HD11	1:A:669:GLU:HG3	1.90	0.52
1:F:329:THR:O	1:F:333:ARG:HG3	2.09	0.52
1:A:661:GLY:N	1:A:664:ASP:OD2	2.39	0.52
1:B:102:GLY:O	1:B:368:SER:HB2	2.09	0.52
1:F:229:MET:HG3	1:F:262:VAL:HG13	1.91	0.52
1:E:517:GLY:HA3	1:E:557:VAL:O	2.09	0.52
1:B:77:PRO:HD2	1:B:337:PHE:CZ	2.45	0.52
1:C:234:SER:HA	1:C:239:PRO:HA	1.92	0.52
1:D:18:GLU:OE2	1:D:325:ARG:NE	2.34	0.52
1:D:5:PRO:HG2	1:D:19:ASP:OD1	2.10	0.52
1:E:137:CYS:SG	1:E:186:HIS:HB2	2.49	0.52
1:A:95:THR:HA	1:A:582:MET:CE	2.40	0.51
1:E:270:ARG:HD3	1:E:274:GLU:OE1	2.11	0.51
1:C:627:ARG:HG2	1:C:628:TYR:O	2.10	0.51
1:A:693:THR:OG1	1:A:695:VAL:O	2.20	0.51
1:A:26:LEU:HD21	1:A:310:VAL:HG11	1.91	0.51
1:A:95:THR:HA	1:A:582:MET:HE1	1.93	0.51
1:B:345:VAL:O	1:B:349:GLN:HG3	2.11	0.51
1:F:43:GLY:HA2	1:F:68:LEU:HD13	1.92	0.51
1:D:5:PRO:HG2	1:D:19:ASP:CG	2.36	0.51
1:D:72:THR:O	1:D:74:LEU:N	2.44	0.51
1:B:100:GLN:HG3	1:B:151:PHE:CD1	2.45	0.51
1:F:367:ARG:NH1	1:F:599:GLU:OE1	2.37	0.51
1:E:408:VAL:HG12	1:E:413:ALA:HB1	1.93	0.51
1:B:204:ILE:HG23	1:B:208:LYS:HD3	1.92	0.51
1:C:111:LEU:HG	1:C:364:VAL:HG22	1.92	0.51
1:D:580:PRO:HG2	1:D:584:GLY:HA3	1.93	0.51
1:B:266:ASP:HA	1:B:299:THR:OG1	2.11	0.51
1:C:266:ASP:HA	1:C:299:THR:OG1	2.11	0.51
1:E:33:MET:HE3	1:E:324:VAL:HG23	1.93	0.51
1:B:576:TRP:CE2	1:B:578:ALA:HB2	2.47	0.50
1:D:290:ARG:HG3	1:D:315:VAL:HG12	1.93	0.50
1:A:202:ALA:HB1	1:A:204:ILE:HD12	1.92	0.50
1:A:232:TYR:HH	2:A:801:BGC:H1	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:VAL:HG12	1:C:165:MET:HG3	1.93	0.50
1:A:660:HIS:NE2	4:A:911:HOH:O	2.35	0.50
1:F:419:GLN:OE1	1:F:444:THR:HA	2.12	0.50
1:E:304:GLU:O	1:E:308:GLU:HG3	2.12	0.50
1:A:576:TRP:CE2	1:A:578:ALA:HB2	2.46	0.50
1:D:3:THR:O	4:D:901:HOH:O	2.20	0.50
1:E:148:ASP:HB3	1:E:559:SER:HB2	1.93	0.50
1:E:264:ASP:OD1	2:E:801:BGC:H1	2.12	0.50
1:F:576:TRP:CZ2	1:F:578:ALA:HB2	2.47	0.50
1:A:137:CYS:SG	1:A:186:HIS:HB2	2.51	0.49
1:E:33:MET:HE2	1:E:323:ALA:HB3	1.94	0.49
1:C:321:ASP:HB3	1:C:325:ARG:HH12	1.76	0.49
1:A:347:ARG:O	1:A:351:VAL:HG22	2.13	0.49
1:C:268:VAL:HA	1:C:271:MET:HE2	1.94	0.49
1:D:347:ARG:O	1:D:351:VAL:HG22	2.12	0.49
1:E:129:HIS:HD2	1:E:341:ARG:HD3	1.76	0.49
1:F:580:PRO:HG2	1:F:584:GLY:HA3	1.95	0.49
1:E:591:LEU:HD11	1:E:598:PRO:HG3	1.92	0.49
1:A:232:TYR:OH	2:A:801:BGC:H1	2.13	0.49
1:A:92:VAL:O	1:A:438:PRO:HD3	2.13	0.49
1:D:137:CYS:HB3	1:D:191:SER:HB3	1.94	0.49
1:A:277:ILE:HA	1:B:693:THR:O	2.11	0.49
1:E:34:LEU:HD21	1:E:302:PHE:CD2	2.48	0.49
1:B:234:SER:HA	1:B:239:PRO:HA	1.94	0.49
1:E:336:LEU:HA	1:E:339:ASP:O	2.12	0.49
1:C:69:THR:O	1:C:75:ARG:HG2	2.13	0.49
1:F:742:GLU:HG2	1:F:743:PHE:N	2.28	0.49
1:A:53:SER:CB	1:A:295:MET:HE1	2.41	0.48
1:C:100:GLN:HG2	1:C:151:PHE:CE1	2.48	0.48
1:D:576:TRP:CZ2	1:D:578:ALA:HB2	2.47	0.48
1:E:100:GLN:HG3	1:E:151:PHE:CD1	2.48	0.48
1:E:669:GLU:HG2	1:E:719:GLY:HA2	1.93	0.48
1:C:425:GLY:HA2	1:C:433:LEU:HD11	1.95	0.48
1:D:102:GLY:O	1:D:368:SER:HB2	2.14	0.48
1:A:409:VAL:HG21	1:A:545:LEU:HD11	1.96	0.48
1:B:408:VAL:HG12	1:B:413:ALA:HB1	1.95	0.48
1:E:57:THR:HG23	1:E:61:ASN:HB2	1.96	0.48
1:C:137:CYS:HB3	1:C:191:SER:HB3	1.96	0.48
1:C:301:ARG:HG2	1:C:304:GLU:OE1	2.13	0.48
1:F:645:TYR:HB3	1:F:678:ARG:HD2	1.96	0.48
1:E:561:PRO:HB3	1:E:606:PHE:CD1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:CYS:HB3	1:A:191:SER:HB3	1.96	0.48
1:B:92:VAL:O	1:B:438:PRO:HD3	2.13	0.48
1:A:234:SER:HA	1:A:239:PRO:HA	1.95	0.48
1:C:88:HIS:CD2	1:C:123:VAL:HG22	2.49	0.48
1:F:576:TRP:CE2	1:F:578:ALA:HB2	2.48	0.48
1:A:229:MET:HE2	1:A:297:MET:HE1	1.96	0.48
1:E:360:VAL:O	1:E:364:VAL:HG23	2.14	0.48
1:E:312:ARG:HB2	1:E:314:LEU:HG	1.94	0.47
1:E:398:PRO:HG2	1:E:401:ALA:HB3	1.95	0.47
1:E:576:TRP:CE2	1:E:578:ALA:HB2	2.49	0.47
1:B:6:TYR:HA	1:B:16:ARG:HB3	1.96	0.47
1:D:19:ASP:O	1:D:23:ARG:HG3	2.15	0.47
1:A:425:GLY:HA2	1:A:433:LEU:HD12	1.96	0.47
1:C:742:GLU:OE2	1:C:762:THR:OG1	2.28	0.47
1:E:92:VAL:O	1:E:438:PRO:HD3	2.14	0.47
1:F:304:GLU:H	1:F:304:GLU:CD	2.22	0.47
1:F:63:LEU:HB3	4:F:1137:HOH:O	2.14	0.46
1:B:73:ARG:CZ	1:B:74:LEU:HD11	2.45	0.46
1:A:316:GLU:HG3	1:A:318:ALA:H	1.80	0.46
1:E:27:PRO:HB3	1:E:307:LEU:HD11	1.97	0.46
1:B:218:GLU:HB2	1:B:255:TRP:CZ2	2.50	0.46
1:E:88:HIS:CD2	1:E:91:TRP:HB2	2.50	0.46
1:C:297:MET:HE1	2:C:801:BGC:C6	2.46	0.46
1:A:746:ARG:NH2	1:A:753:GLU:OE2	2.47	0.46
1:C:696:THR:OG1	1:F:237:GLY:HA2	2.16	0.46
1:F:325:ARG:HD2	4:F:998:HOH:O	2.16	0.45
1:E:102:GLY:O	1:E:368:SER:HB2	2.16	0.45
1:E:176:ASP:O	1:E:341:ARG:NH2	2.49	0.45
1:A:732:ALA:C	1:A:734:GLY:H	2.24	0.45
1:D:163:SER:OG	1:D:167:ARG:NH1	2.49	0.45
1:B:20:LEU:O	1:B:24:MET:HG3	2.17	0.45
1:E:347:ARG:O	1:E:351:VAL:HG22	2.16	0.45
1:D:561:PRO:HB3	1:D:606:PHE:CD1	2.51	0.45
1:B:8:ASP:HB3	1:B:11:VAL:HG23	1.99	0.45
1:A:450:ARG:NH1	1:A:459:VAL:O	2.50	0.45
1:C:329:THR:CG2	1:C:333:ARG:HE	2.24	0.45
1:D:264:ASP:OD1	2:D:801:BGC:H1	2.17	0.45
1:B:333:ARG:NH1	4:B:924:HOH:O	2.46	0.45
1:C:112:VAL:CG1	1:C:165:MET:HG3	2.47	0.45
1:F:408:VAL:HG12	1:F:413:ALA:HB1	1.98	0.45
1:E:576:TRP:CZ2	1:E:578:ALA:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:GLU:OE1	1:C:325:ARG:NE	2.46	0.44
1:E:356:GLU:H	1:E:356:GLU:HG2	1.46	0.44
1:A:5:PRO:HG2	1:A:19:ASP:OD1	2.17	0.44
1:A:638:ALA:O	1:A:641:GLU:HG2	2.17	0.44
1:B:450:ARG:NH1	1:B:459:VAL:O	2.49	0.44
1:D:387:ALA:N	1:D:396:LEU:HD23	2.32	0.44
1:C:151:PHE:HB2	1:C:158:ILE:HG12	1.98	0.44
1:D:232:TYR:OH	1:D:265:TRP:HB3	2.17	0.44
1:D:576:TRP:CE2	1:D:578:ALA:HB2	2.53	0.44
1:B:645:TYR:HB3	1:B:678:ARG:HD2	1.99	0.44
1:B:731:ASP:OD1	1:B:732:ALA:O	2.36	0.44
1:E:54:LEU:HG	1:E:78:LEU:HD12	1.98	0.44
1:A:119:THR:O	1:A:123:VAL:HG23	2.17	0.44
1:A:189:GLY:HA3	1:A:235:MET:HE1	1.99	0.44
1:C:290:ARG:NH2	4:C:915:HOH:O	2.38	0.44
1:B:34:LEU:HB3	1:B:36:LEU:HD11	2.00	0.44
1:C:423:TRP:CZ3	1:C:429:GLN:HG2	2.52	0.44
1:F:623:GLN:HG3	1:F:625:GLY:O	2.18	0.44
1:E:310:VAL:HG22	1:E:315:VAL:HG23	2.00	0.44
1:A:21:LEU:HD11	1:A:29:LYS:HD3	1.99	0.43
1:B:26:LEU:HD21	1:B:310:VAL:CG2	2.37	0.43
1:B:187:PHE:O	1:B:240:VAL:HG11	2.18	0.43
1:E:76:ILE:HG23	1:E:337:PHE:CZ	2.53	0.43
1:E:268:VAL:HA	1:E:271:MET:HE2	1.99	0.43
1:B:257:TYR:CZ	1:B:259:GLY:HA3	2.53	0.43
1:E:265:TRP:O	1:E:270:ARG:NH1	2.52	0.43
1:F:171:GLY:C	1:F:173:GLY:H	2.27	0.43
1:B:306:ALA:O	1:B:310:VAL:HG22	2.18	0.43
1:A:90:PHE:HA	1:A:433:LEU:HD11	2.01	0.43
1:C:398:PRO:HG2	1:C:401:ALA:HB3	2.00	0.43
1:E:170:GLN:HG2	1:E:180:ILE:O	2.19	0.43
1:C:273:TRP:CZ3	1:C:274:GLU:HB2	2.54	0.43
1:E:138:ILE:HD11	1:E:153:GLU:O	2.19	0.43
1:A:148:ASP:HB3	1:A:559:SER:HB2	2.01	0.42
1:B:20:LEU:HD11	1:B:74:LEU:HD23	2.01	0.42
1:B:347:ARG:HE	1:B:351:VAL:HG11	1.84	0.42
1:C:578:ALA:O	1:C:579:ASN:C	2.62	0.42
1:F:138:ILE:HD11	1:F:153:GLU:O	2.19	0.42
1:E:119:THR:O	1:E:123:VAL:HG23	2.19	0.42
1:F:102:GLY:O	1:F:368:SER:HB2	2.20	0.42
1:D:301:ARG:HG2	1:D:301:ARG:NH1	2.26	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:656:LEU:HD11	1:B:669:GLU:HG3	2.01	0.42
1:F:507:ARG:HD3	4:F:1264:HOH:O	2.19	0.42
1:E:301:ARG:HA	1:E:304:GLU:OE2	2.18	0.42
1:E:20:LEU:HD11	1:E:74:LEU:HD13	2.01	0.42
1:B:372:LEU:HA	1:B:372:LEU:HD23	1.83	0.42
1:B:732:ALA:C	1:B:734:GLY:H	2.28	0.42
1:B:511:LEU:HD11	1:B:553:VAL:HG23	2.01	0.42
1:F:171:GLY:O	4:F:901:HOH:O	2.21	0.42
1:E:33:MET:CE	1:E:323:ALA:HB3	2.50	0.42
1:A:398:PRO:O	4:A:901:HOH:O	2.21	0.42
1:B:4:LEU:HD22	1:B:23:ARG:HH12	1.84	0.42
1:B:137:CYS:SG	1:B:186:HIS:HB2	2.60	0.42
1:E:8:ASP:C	1:E:10:ALA:H	2.28	0.42
1:E:321:ASP:HB3	1:E:325:ARG:NH2	2.35	0.42
1:B:204:ILE:HD12	1:B:235:MET:HE2	2.02	0.41
1:C:174:LEU:HD13	1:C:333:ARG:HB2	2.01	0.41
1:F:693:THR:OG1	1:F:695:VAL:O	2.32	0.41
1:D:423:TRP:CH2	1:D:429:GLN:HG2	2.55	0.41
1:E:746:ARG:HA	1:E:757:LEU:O	2.20	0.41
1:E:753:GLU:HA	1:E:756:LEU:HD12	2.02	0.41
1:C:5:PRO:O	1:C:16:ARG:HG2	2.20	0.41
1:C:148:ASP:HB3	1:C:559:SER:HB2	2.02	0.41
1:A:200:SER:HB2	4:A:1142:HOH:O	2.20	0.41
1:E:245:TRP:O	1:E:249:ASP:HB2	2.20	0.41
1:E:623:GLN:HG2	1:E:625:GLY:O	2.20	0.41
1:A:304:GLU:H	1:A:304:GLU:CD	2.28	0.41
1:B:247:LEU:O	1:B:252:ARG:HG2	2.21	0.41
1:C:325:ARG:HB2	1:C:326:ARG:NH1	2.34	0.41
1:D:148:ASP:HB3	1:D:559:SER:HB2	2.01	0.41
1:E:30:VAL:HG12	1:E:303:PHE:HD1	1.85	0.41
1:C:100:GLN:HG2	1:C:151:PHE:CD1	2.56	0.41
1:B:26:LEU:HD13	1:B:317:GLU:CD	2.46	0.41
1:B:745:LEU:O	1:B:758:ARG:HA	2.20	0.41
1:F:333:ARG:HH21	1:F:333:ARG:HD2	1.76	0.41
1:D:132:PHE:HB3	1:D:185:LYS:HE3	2.03	0.41
1:E:36:LEU:HD12	1:E:51:ALA:HB2	2.02	0.41
1:E:57:THR:CG2	1:E:61:ASN:HB2	2.50	0.41
1:C:326:ARG:N	1:C:326:ARG:HD2	2.36	0.41
1:F:732:ALA:O	1:F:733:HIS:CG	2.73	0.41
1:B:14:ALA:O	1:B:18:GLU:HG2	2.21	0.41
1:B:20:LEU:CD1	1:B:74:LEU:HD23	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:ILE:HG12	1:B:582:MET:HE3	2.02	0.41
1:B:261:LEU:O	1:B:294:ASP:HB2	2.21	0.41
1:C:732:ALA:O	1:C:733:HIS:HB2	2.21	0.41
1:F:58:SER:HB2	1:F:60:GLU:OE2	2.21	0.41
1:D:60:GLU:CD	1:D:60:GLU:H	2.29	0.41
1:D:387:ALA:HB2	1:D:396:LEU:HD21	2.02	0.41
1:D:627:ARG:HA	1:D:633:GLN:HG3	2.03	0.41
1:E:56:HIS:HA	1:E:82:GLU:OE1	2.20	0.41
1:E:321:ASP:HB3	1:E:325:ARG:HH21	1.86	0.41
1:A:214:LEU:HA	1:A:214:LEU:HD23	1.84	0.41
1:F:584:GLY:O	1:F:588:VAL:HG23	2.21	0.41
1:D:92:VAL:O	1:D:438:PRO:HD3	2.21	0.41
1:F:270:ARG:NH2	1:F:274:GLU:OE1	2.54	0.40
1:D:409:VAL:HG12	1:D:462:ALA:HB3	2.04	0.40
1:B:31:GLY:O	1:B:34:LEU:HB2	2.21	0.40
1:C:31:GLY:HA2	1:C:34:LEU:HD12	2.03	0.40
1:F:4:LEU:HD23	1:F:7:LEU:HD23	2.03	0.40
1:B:248:ASP:O	1:B:253:GLY:N	2.48	0.40
1:D:306:ALA:O	1:D:310:VAL:HG23	2.21	0.40
1:A:42:VAL:HG12	1:A:68:LEU:CD1	2.50	0.40
1:B:55:LEU:HA	1:B:81:ALA:O	2.21	0.40
1:F:616:TYR:O	1:F:619:VAL:HG23	2.20	0.40
1:A:614:PRO:HD2	1:B:197:ARG:CZ	2.51	0.40
1:A:616:TYR:O	1:A:619:VAL:HG23	2.20	0.40
1:B:151:PHE:HB2	1:B:158:ILE:HG12	2.03	0.40
1:C:433:LEU:HG	1:C:433:LEU:O	2.22	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	743/768 (97%)	711 (96%)	31 (4%)	1 (0%)	48	64
1	B	740/768 (96%)	703 (95%)	36 (5%)	1 (0%)	48	64
1	C	741/768 (96%)	700 (94%)	38 (5%)	3 (0%)	30	43
1	D	742/768 (97%)	705 (95%)	32 (4%)	5 (1%)	18	28
1	E	728/768 (95%)	687 (94%)	36 (5%)	5 (1%)	18	28
1	F	739/768 (96%)	714 (97%)	15 (2%)	10 (1%)	9	13
All	All	4433/4608 (96%)	4220 (95%)	188 (4%)	25 (1%)	21	32

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	186	HIS
1	F	416	ASP
1	D	73	ARG
1	D	416	ASP
1	E	270	ARG
1	B	186	HIS
1	F	4	LEU
1	F	73	ARG
1	F	392	ASP
1	E	92	VAL
1	A	186	HIS
1	F	40	ASP
1	F	72	THR
1	F	172	ASP
1	F	186	HIS
1	F	733	HIS
1	D	4	LEU
1	E	416	ASP
1	E	733	HIS
1	C	657	GLY
1	C	733	HIS
1	F	303	PHE
1	D	40	ASP
1	E	56	HIS
1	D	474	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	574/589 (98%)	572 (100%)	2 (0%)	86	93
1	B	570/589 (97%)	568 (100%)	2 (0%)	84	92
1	C	573/589 (97%)	572 (100%)	1 (0%)	87	94
1	D	569/589 (97%)	569 (100%)	0	100	100
1	E	564/589 (96%)	559 (99%)	5 (1%)	70	85
1	F	569/589 (97%)	566 (100%)	3 (0%)	81	91
All	All	3419/3534 (97%)	3406 (100%)	13 (0%)	84	92

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	433	LEU
1	A	435	ASP
1	B	92	VAL
1	B	652	ASP
1	C	692	VAL
1	F	78	LEU
1	F	92	VAL
1	F	473	ASP
1	E	16	ARG
1	E	67	GLU
1	E	92	VAL
1	E	475	GLU
1	E	720	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	349	GLN
1	B	633	GLN
1	C	186	HIS
1	F	117	HIS

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Mol	Chain	Res	Type
1	F	276	HIS
1	F	278	GLN
1	F	293	ASN
1	F	348	GLN
1	F	349	GLN
1	D	88	HIS
1	D	293	ASN
1	D	487	GLN
1	E	35	GLN
1	E	61	ASN
1	E	186	HIS
1	E	275	GLN
1	E	623	GLN
1	E	633	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](#)

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BGC	E	801	-	12,12,12	1.22	0	17,17,17	0.78	0
2	BGC	B	801	-	12,12,12	1.18	1 (8%)	17,17,17	1.93	4 (23%)
2	BGC	F	801	-	12,12,12	1.20	0	17,17,17	1.36	3 (17%)
2	BGC	A	801	-	12,12,12	1.46	3 (25%)	17,17,17	1.61	5 (29%)
2	BGC	D	801	-	12,12,12	1.18	0	17,17,17	0.62	0
2	BGC	C	801	-	12,12,12	1.29	1 (8%)	17,17,17	1.64	2 (11%)

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	E	801	-	-	2/2/22/22	0/1/1/1
2	BGC	B	801	-	-	2/2/22/22	0/1/1/1
2	BGC	F	801	-	-	2/2/22/22	0/1/1/1
2	BGC	A	801	-	-	0/2/22/22	0/1/1/1
2	BGC	D	801	-	-	2/2/22/22	0/1/1/1
2	BGC	C	801	-	-	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	A	801	BGC	O5-C5	2.31	1.50	1.44
2	B	801	BGC	C3-C2	-2.29	1.46	1.52
2	A	801	BGC	C1-C2	2.25	1.57	1.52
2	A	801	BGC	O5-C1	2.08	1.48	1.42
2	C	801	BGC	C3-C2	-2.01	1.47	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	BGC	C1-O5-C5	-4.62	104.72	113.65
2	C	801	BGC	O5-C1-C2	4.39	118.02	110.30
2	B	801	BGC	O5-C1-C2	4.08	117.47	110.30
2	B	801	BGC	O3-C3-C2	-3.42	102.32	110.38
2	A	801	BGC	O5-C1-C2	3.11	115.76	110.30
2	A	801	BGC	C1-O5-C5	2.82	119.11	113.65
2	F	801	BGC	O5-C1-C2	2.77	115.17	110.30
2	A	801	BGC	O5-C5-C6	2.75	113.25	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	801	BGC	O5-C5-C4	2.60	114.38	109.70
2	F	801	BGC	O3-C3-C2	-2.30	104.95	110.38
2	A	801	BGC	O5-C5-C4	2.25	113.75	109.70
2	F	801	BGC	C1-O5-C5	-2.25	109.30	113.65
2	B	801	BGC	O1-C1-C2	2.12	115.14	108.98
2	A	801	BGC	O1-C1-C2	2.02	114.86	108.98

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	801	BGC	O5-C5-C6-O6
2	C	801	BGC	C4-C5-C6-O6
2	E	801	BGC	C4-C5-C6-O6
2	B	801	BGC	O5-C5-C6-O6
2	B	801	BGC	C4-C5-C6-O6
2	E	801	BGC	O5-C5-C6-O6
2	F	801	BGC	C4-C5-C6-O6
2	F	801	BGC	O5-C5-C6-O6
2	D	801	BGC	O5-C5-C6-O6
2	D	801	BGC	C4-C5-C6-O6

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	801	BGC	2	0
2	B	801	BGC	1	0
2	A	801	BGC	2	0
2	D	801	BGC	1	0
2	C	801	BGC	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	749/768 (97%)	-0.41	10 (1%) 75 71	24, 36, 56, 99	0
1	B	746/768 (97%)	-0.14	19 (2%) 58 54	25, 39, 79, 95	0
1	C	747/768 (97%)	-0.25	25 (3%) 49 45	23, 35, 74, 94	0
1	D	748/768 (97%)	-0.33	18 (2%) 59 55	28, 39, 60, 100	0
1	E	740/768 (96%)	0.23	86 (11%) 9 7	28, 45, 108, 125	0
1	F	747/768 (97%)	-0.47	11 (1%) 72 68	23, 34, 54, 89	0
All	All	4477/4608 (97%)	-0.23	169 (3%) 44 40	23, 38, 78, 125	0

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	392	ASP	6.8
1	E	173	GLY	5.8
1	E	42	VAL	5.0
1	E	46	VAL	5.0
1	B	387	ALA	4.9
1	D	387	ALA	4.9
1	F	765	GLY	4.8
1	E	474	PRO	4.8
1	E	92	VAL	4.8
1	E	6	TYR	4.7
1	E	38	ALA	4.7
1	B	473	ASP	4.5
1	F	175	SER	4.4
1	E	319	ALA	4.4
1	D	474	PRO	4.4
1	E	473	ASP	4.3
1	E	307	LEU	4.3
1	C	391	PRO	4.3
1	A	764	ALA	4.3

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Mol	Chain	Res	Type	RSRZ
1	E	12	PRO	4.1
1	E	310	VAL	4.0
1	E	7	LEU	3.9
1	A	392	ASP	3.9
1	E	63	LEU	3.8
1	D	391	PRO	3.8
1	E	58	SER	3.7
1	E	175	SER	3.7
1	E	19	ASP	3.7
1	E	43	GLY	3.7
1	E	26	LEU	3.7
1	E	74	LEU	3.6
1	F	391	PRO	3.6
1	E	76	ILE	3.5
1	E	303	PHE	3.5
1	D	172	ASP	3.5
1	F	388	ALA	3.4
1	D	175	SER	3.4
1	A	391	PRO	3.4
1	B	10	ALA	3.4
1	A	659	GLU	3.4
1	B	3	THR	3.4
1	E	41	GLY	3.4
1	F	93	GLY	3.3
1	F	473	ASP	3.3
1	B	12	PRO	3.3
1	F	173	GLY	3.3
1	E	44	PRO	3.3
1	E	320	ILE	3.2
1	E	311	ASP	3.2
1	A	387	ALA	3.2
1	C	5	PRO	3.2
1	E	314	LEU	3.2
1	E	302	PHE	3.1
1	A	652	ASP	3.1
1	E	337	PHE	3.1
1	E	33	MET	3.1
1	E	300	PRO	3.1
1	E	765	GLY	3.1
1	C	11	VAL	3.1
1	E	281	TYR	3.0
1	E	59	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	332	PHE	3.0
1	E	13	VAL	3.0
1	C	387	ALA	3.0
1	E	282	VAL	3.0
1	E	77	PRO	3.0
1	E	49	LYS	3.0
1	C	30	VAL	2.9
1	C	316	GLU	2.9
1	E	68	LEU	2.9
1	B	765	GLY	2.9
1	E	73	ARG	2.9
1	B	317	GLU	2.9
1	E	393	GLY	2.9
1	C	12	PRO	2.8
1	E	72	THR	2.8
1	B	92	VAL	2.8
1	C	92	VAL	2.8
1	D	3	THR	2.8
1	F	472	PRO	2.8
1	E	30	VAL	2.8
1	B	14	ALA	2.8
1	E	78	LEU	2.8
1	E	48	GLU	2.8
1	C	93	GLY	2.8
1	E	22	ALA	2.8
1	B	472	PRO	2.7
1	E	57	THR	2.7
1	E	299	THR	2.7
1	E	475	GLU	2.7
1	E	315	VAL	2.7
1	E	27	PRO	2.7
1	A	92	VAL	2.7
1	E	11	VAL	2.7
1	E	306	ALA	2.7
1	C	317	GLU	2.7
1	D	487	GLN	2.7
1	C	392	ASP	2.6
1	B	11	VAL	2.6
1	E	24	MET	2.6
1	A	474	PRO	2.6
1	D	473	ASP	2.6
1	E	280	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	13	VAL	2.6
1	E	272	VAL	2.6
1	E	51	ALA	2.6
1	D	173	GLY	2.5
1	D	393	GLY	2.5
1	E	37	ASP	2.5
1	E	60	GLU	2.5
1	E	14	ALA	2.5
1	C	313	GLY	2.5
1	D	174	LEU	2.5
1	D	764	ALA	2.5
1	E	8	ASP	2.5
1	E	15	ASP	2.5
1	C	321	ASP	2.4
1	F	392	ASP	2.4
1	E	23	ARG	2.4
1	A	658	THR	2.4
1	E	21	LEU	2.4
1	F	432	TRP	2.4
1	E	308	GLU	2.4
1	C	14	ALA	2.4
1	C	475	GLU	2.4
1	E	71	ARG	2.4
1	E	64	ALA	2.4
1	E	305	GLY	2.3
1	B	393	GLY	2.3
1	E	334	LEU	2.3
1	D	416	ASP	2.3
1	F	92	VAL	2.3
1	E	312	ARG	2.3
1	D	41	GLY	2.3
1	B	316	GLU	2.3
1	C	13	VAL	2.3
1	C	393	GLY	2.2
1	C	20	LEU	2.2
1	E	304	GLU	2.2
1	C	474	PRO	2.2
1	C	22	ALA	2.2
1	D	433	LEU	2.2
1	C	332	PHE	2.2
1	D	392	ASP	2.2
1	B	22	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	322	ALA	2.1
1	D	578	ALA	2.1
1	E	50	HIS	2.1
1	C	174	LEU	2.1
1	E	328	LEU	2.1
1	C	652	ASP	2.1
1	E	268	VAL	2.1
1	C	651	ALA	2.1
1	E	69	THR	2.1
1	D	430	ALA	2.1
1	E	32	GLN	2.1
1	E	324	VAL	2.1
1	E	34	LEU	2.1
1	B	93	GLY	2.1
1	E	274	GLU	2.1
1	C	71	ARG	2.0
1	E	36	LEU	2.0
1	E	338	GLU	2.0
1	E	75	ARG	2.0
1	E	386	ARG	2.0
1	B	47	LEU	2.0
1	E	29	LYS	2.0
1	A	3	THR	2.0
1	E	329	THR	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](#)

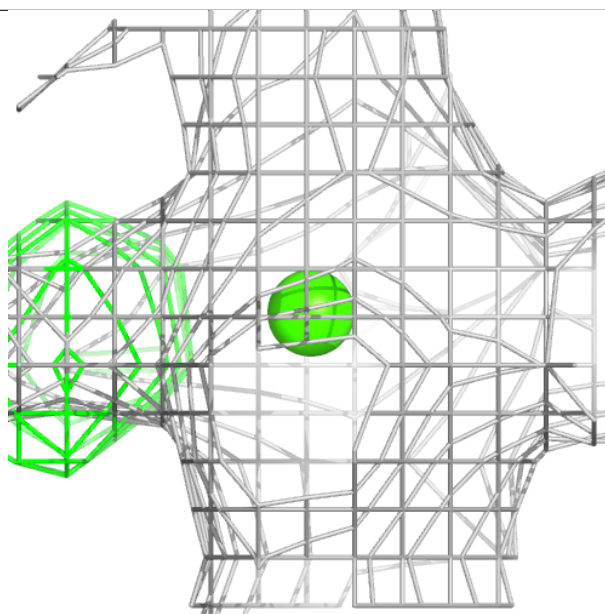
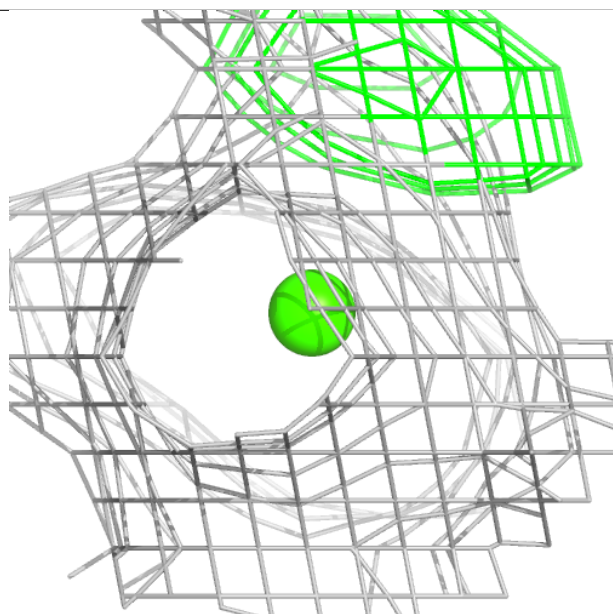
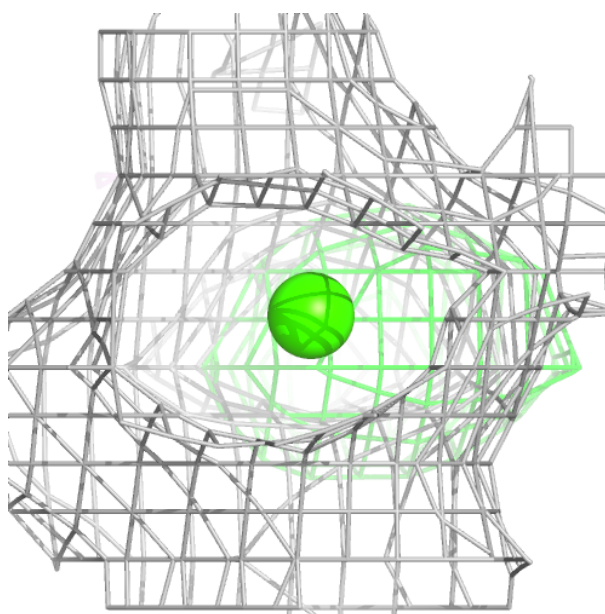
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B	801	12/12	0.80	0.14	38,45,50,56	0
2	BGC	E	801	12/12	0.91	0.08	45,51,56,59	0
2	BGC	F	801	12/12	0.92	0.09	28,36,39,39	0
2	BGC	C	801	12/12	0.93	0.11	30,39,46,47	0
2	BGC	A	801	12/12	0.94	0.13	34,35,43,45	0
2	BGC	D	801	12/12	0.96	0.06	34,40,46,47	0
3	CA	C	802	1/1	0.96	0.05	50,50,50,50	0
3	CA	B	802	1/1	0.98	0.05	48,48,48,48	0
3	CA	F	802	1/1	0.98	0.04	52,52,52,52	0
3	CA	A	802	1/1	0.99	0.04	60,60,60,60	0
3	CA	D	802	1/1	0.99	0.04	58,58,58,58	0
3	CA	E	802	1/1	0.99	0.05	58,58,58,58	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.

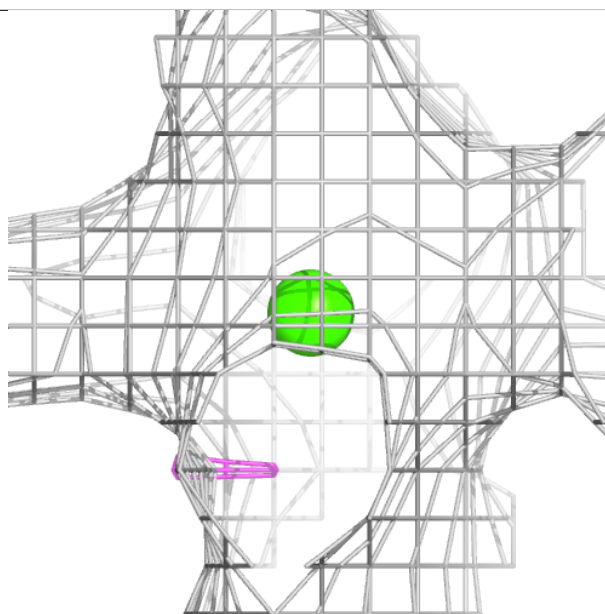
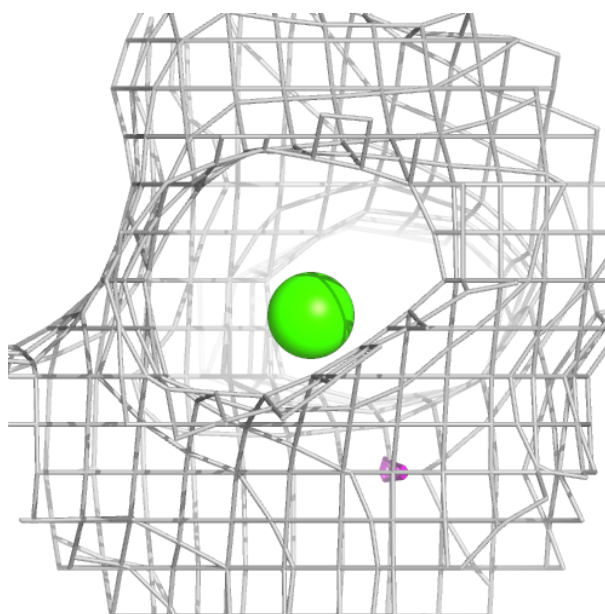
Electron density around CA C 802:

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



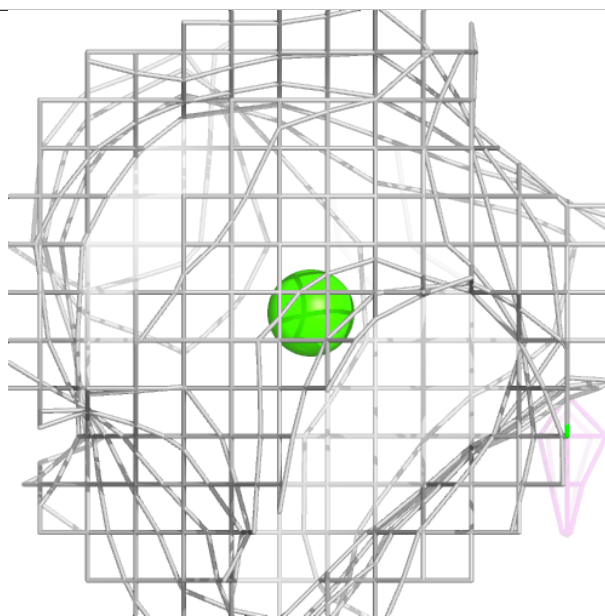
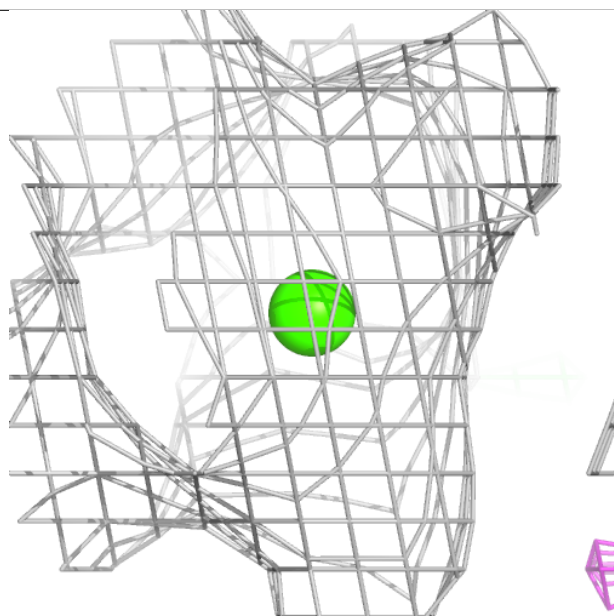
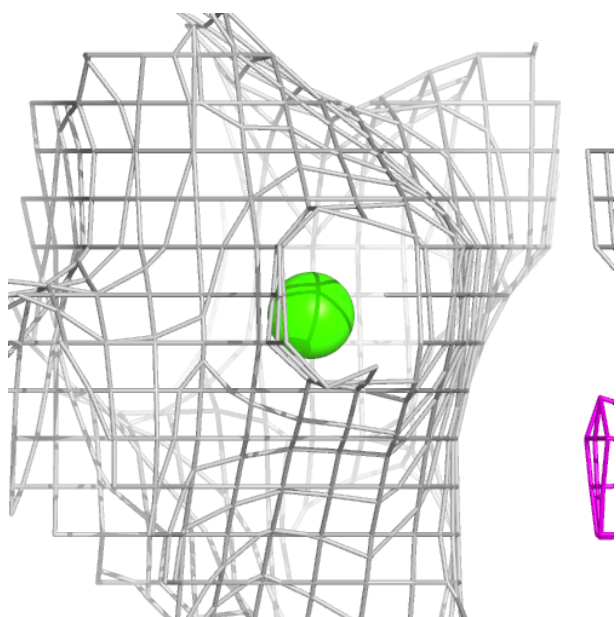
Electron density around CA B 802:

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



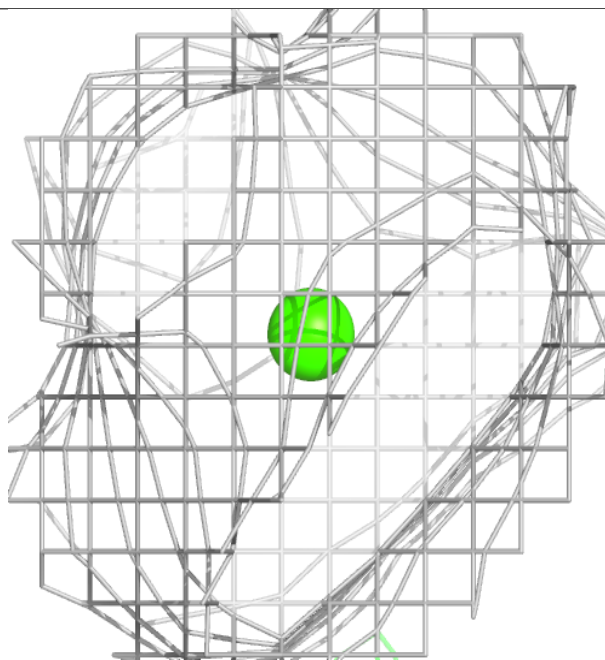
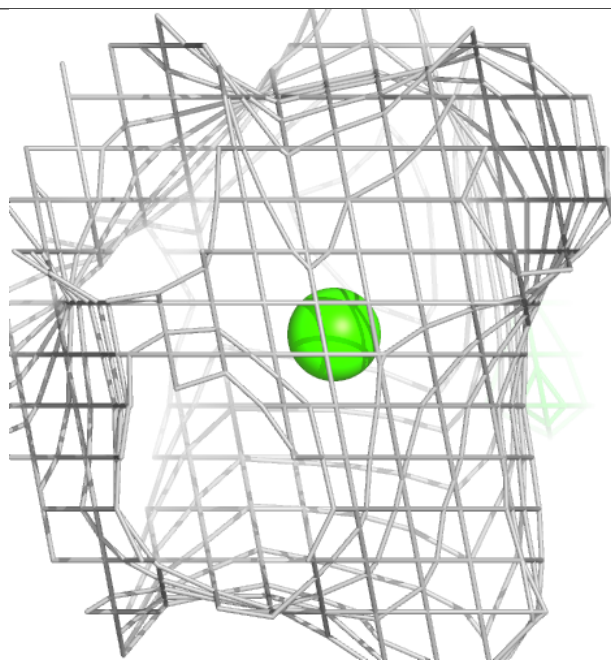
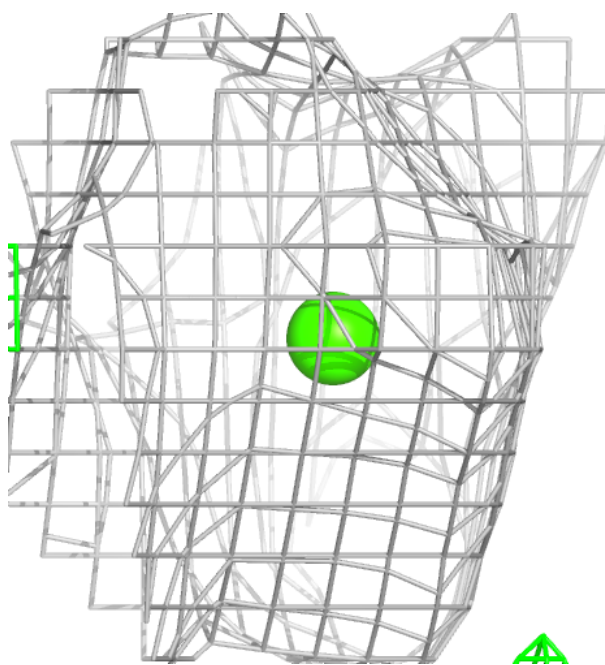
Electron density around CA F 802:

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



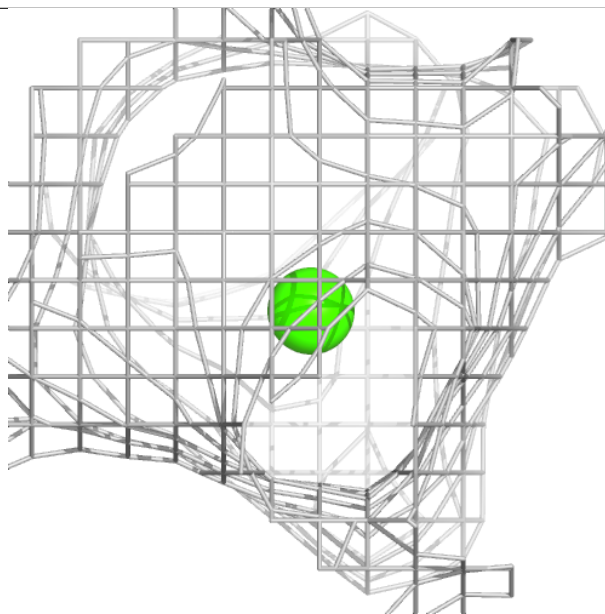
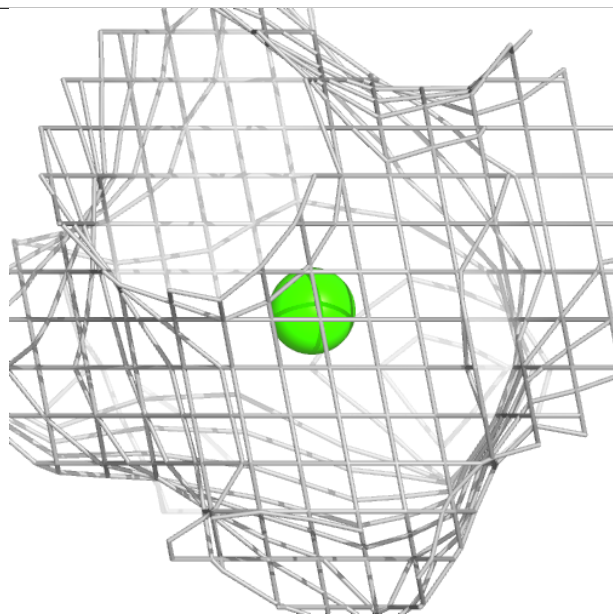
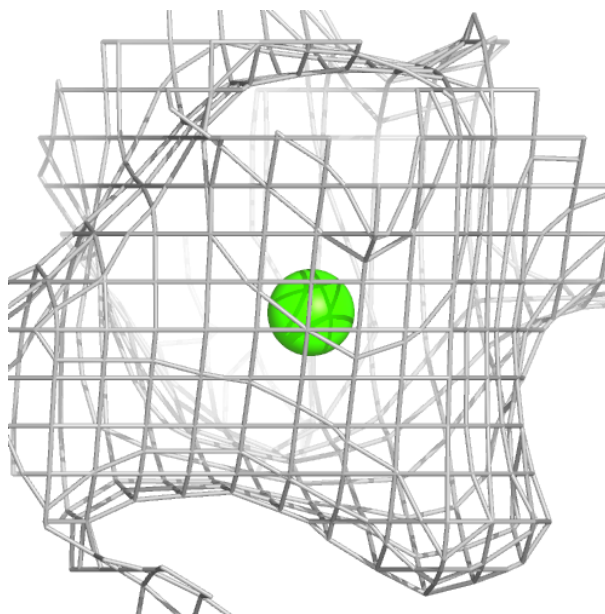
Electron density around CA A 802:

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



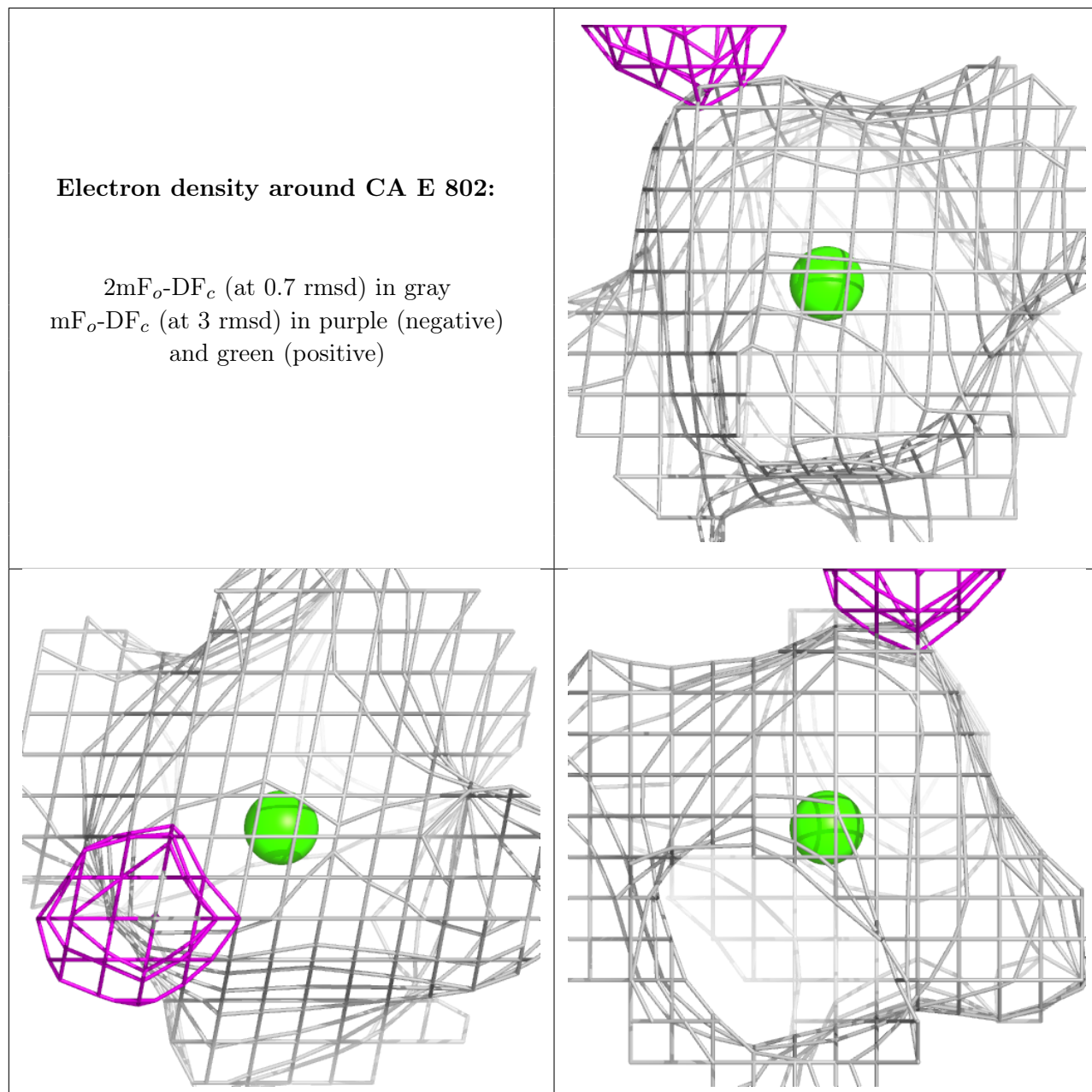
Electron density around CA D 802:

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



Electron density around CA E 802:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.