



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 14, 2025 – 01:33 pm BST

PDB ID : 8QRZ / pdb_00008qrz
Title : Microbacterium testaceum C-glucosyl deglycosidase (CGD), wild type
Authors : Furlanetto, V.; Kalyani, D.C.; Divne, C.
Deposited on : 2023-10-09
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

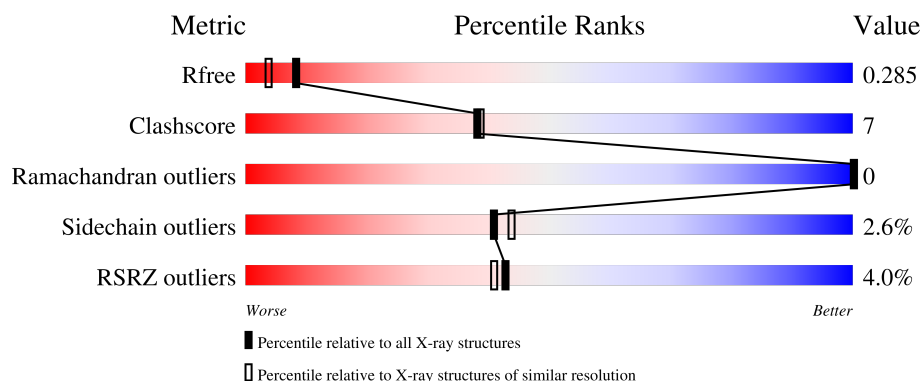
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	435	<div> <div>0%</div> <div>85%</div> <div>15%</div> <div>.</div> </div>
1	B	435	<div> <div>6%</div> <div>87%</div> <div>13%</div> <div>.</div> </div>
1	C	435	<div> <div>3%</div> <div>81%</div> <div>17%</div> <div>.</div> </div>
1	D	435	<div> <div>6%</div> <div>79%</div> <div>20%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	RB	A	501	-	-	X	-
2	RB	B	501	-	-	X	-
2	RB	C	501	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14165 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 Sugar phosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	435	Total	C	N	O	S	0	1	0
			3451	2203	592	646	10			
1	B	435	Total	C	N	O	S	0	1	0
			3448	2202	591	645	10			
1	C	435	Total	C	N	O	S	0	1	0
			3448	2202	591	645	10			
1	D	435	Total	C	N	O	S	0	0	0
			3440	2197	588	645	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	434	ALA	-	expression tag	UNP A0A147EV52
A	435	GLU	-	expression tag	UNP A0A147EV52
B	434	ALA	-	expression tag	UNP A0A147EV52
B	435	GLU	-	expression tag	UNP A0A147EV52
C	434	ALA	-	expression tag	UNP A0A147EV52
C	435	GLU	-	expression tag	UNP A0A147EV52
D	434	ALA	-	expression tag	UNP A0A147EV52
D	435	GLU	-	expression tag	UNP A0A147EV52

- Molecule 2 is RUBIDIUM ION (CCD ID: RB) (formula: Rb) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Rb	0	0
			1	1		
2	B	1	Total	Rb	0	0
			1	1		
2	C	1	Total	Rb	0	0
			1	1		
2	D	1	Total	Rb	0	0
			1	1		

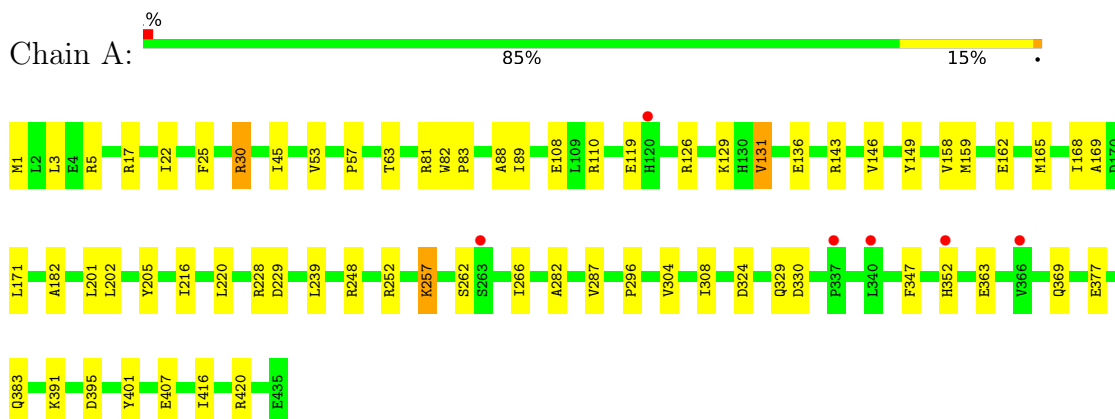
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	99	Total 99	O 99	0	0
3	B	104	Total 104	O 104	0	0
3	C	84	Total 84	O 84	0	0
3	D	87	Total 87	O 87	0	0

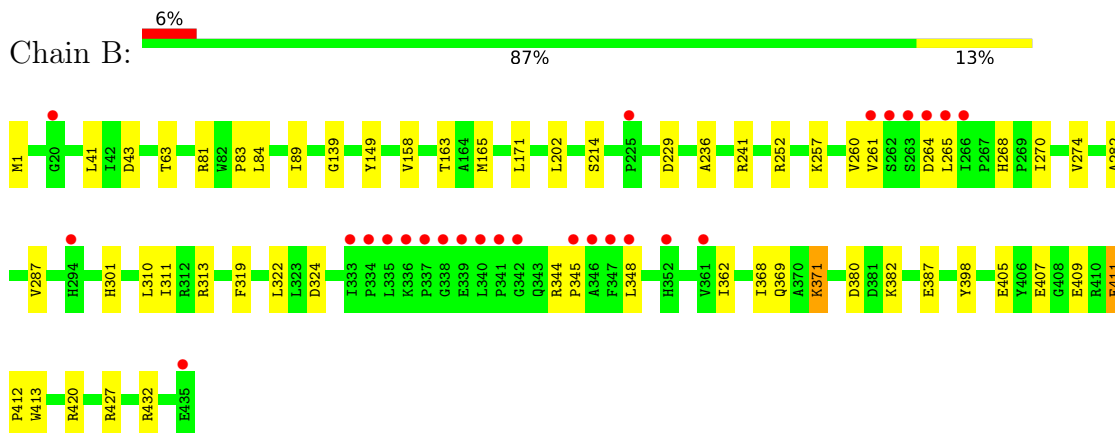
3 Residue-property plots [i](#)

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.

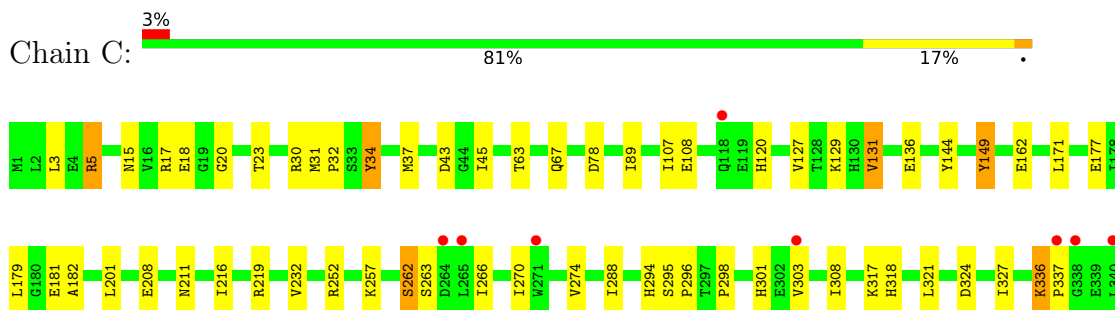
- Molecule 1: Sugar phosphate isomerase

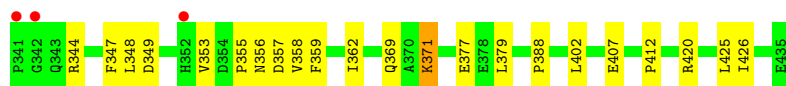


- Molecule 1: Sugar phosphate isomerase

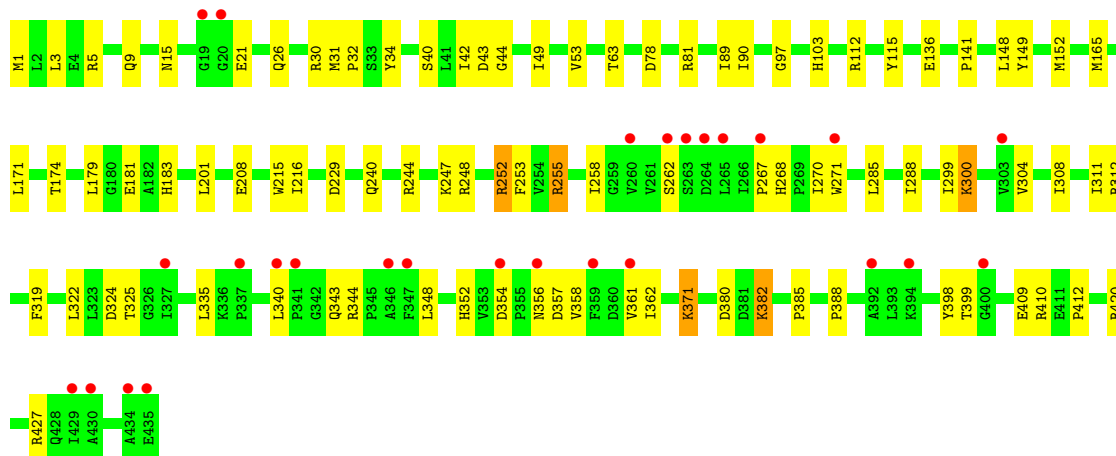
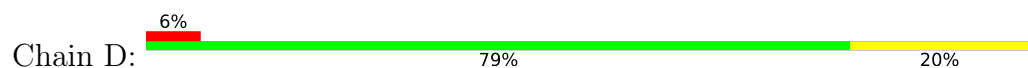


- Molecule 1: Sugar phosphate isomerase





• Molecule 1: Sugar phosphate isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	78.22Å 80.41Å 82.96Å 90.17° 94.39° 106.58°	Depositor
Resolution (Å)	39.91 – 2.00 39.91 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.8 (39.91-2.00) 98.8 (39.91-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.236 , 0.290 0.236 , 0.285	Depositor DCC
R_{free} test set	128473 reflections (1.56%)	wwPDB-VP
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14165	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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:
RB

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	0.43	0/3541	0.65	0/4825
1	B	0.43	0/3541	0.63	0/4825
1	C	0.42	0/3541	0.64	0/4825
1	D	0.40	0/3530	0.65	1/4811 (0.0%)
All	All	0.42	0/14153	0.64	1/19286 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	97	GLY	C-N-CA	6.25	137.31	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3451	0	3396	42	0
1	B	3448	0	3397	37	0
1	C	3448	0	3397	53	0
1	D	3440	0	3384	62	0
2	A	1	0	0	4	0
2	B	1	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	3	0
2	D	1	0	0	1	0
3	A	99	0	0	4	0
3	B	104	0	0	3	0
3	C	84	0	0	6	0
3	D	87	0	0	7	0
All	All	14165	0	13574	186	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 7.

All (186) 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:219:ARG:O	1:D:112:ARG:NH1	2.06	0.88
1:D:40:SER:O	1:D:112:ARG:NH2	2.13	0.82
1:A:5:ARG:O	1:A:30:ARG:NH1	2.16	0.79
1:B:369:GLN:OE1	2:B:501:RB:RB	2.10	0.78
1:A:171:LEU:HD21	1:A:420:ARG:HG2	1.64	0.78
1:A:369:GLN:OE1	2:A:501:RB:RB	2.11	0.76
1:B:165:MET:HE1	1:B:202:LEU:HA	1.69	0.74
1:C:402:LEU:HD12	1:C:426:ILE:HD12	1.70	0.73
1:A:129:LYS:HG2	1:A:131:VAL:HG13	1.70	0.72
1:A:324:ASP:OD2	2:A:501:RB:RB	2.18	0.70
1:D:252:ARG:NH2	3:D:601:HOH:O	2.23	0.69
1:B:324:ASP:OD2	2:B:501:RB:RB	2.19	0.68
1:C:369:GLN:OE1	2:C:501:RB:RB	2.20	0.67
1:B:387:GLU:OE1	1:B:432:ARG:NH2	2.28	0.67
1:C:252:ARG:HG3	3:C:615:HOH:O	1.94	0.66
1:C:324:ASP:OD2	2:C:501:RB:RB	2.22	0.66
1:D:324:ASP:OD2	2:D:501:RB:RB	2.22	0.65
1:B:171:LEU:HD21	1:B:420:ARG:HA	1.79	0.64
1:C:336:LYS:HG2	1:C:337:PRO:HD2	1.79	0.64
1:D:427:ARG:NH1	3:D:604:HOH:O	2.30	0.63
1:C:288:ILE:HG12	1:C:318:HIS:HA	1.80	0.63
1:B:81:ARG:NH1	3:B:603:HOH:O	2.31	0.63
2:A:501:RB:RB	3:A:625:HOH:O	2.25	0.62
1:D:3:LEU:HD13	1:D:30:ARG:HB3	1.81	0.62
1:D:340:LEU:CD1	3:D:602:HOH:O	2.48	0.62
1:A:330:ASP:HA	1:A:352:HIS:CD2	2.34	0.61
1:B:214:SER:OG	3:B:601:HOH:O	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:340:LEU:HD13	3:D:602:HOH:O	2.01	0.61
1:C:371:LYS:NZ	3:C:607:HOH:O	2.33	0.60
1:C:129:LYS:HG2	1:C:131:VAL:HG13	1.84	0.60
1:D:171:LEU:HD21	1:D:420:ARG:HA	1.84	0.60
1:D:247:LYS:HG2	1:D:285:LEU:HD22	1.84	0.59
1:B:322:LEU:HD21	1:B:369:GLN:HG3	1.84	0.59
1:D:49:ILE:H	1:D:53:VAL:HG23	1.66	0.59
2:B:501:RB:RB	3:B:624:HOH:O	2.28	0.59
2:C:501:RB:RB	3:C:662:HOH:O	2.28	0.59
1:C:308:ILE:HD11	1:C:321:LEU:HD11	1.84	0.58
1:D:247:LYS:HE3	1:D:285:LEU:O	2.04	0.58
1:D:362:ILE:HD11	1:D:398:TYR:HB2	1.85	0.58
1:D:255:ARG:HG3	1:D:322:LEU:HD13	1.85	0.58
2:A:501:RB:RB	3:A:639:HOH:O	2.27	0.58
1:D:248:ARG:NH2	3:D:603:HOH:O	2.30	0.57
1:A:248:ARG:NH2	3:A:603:HOH:O	2.31	0.57
1:A:252:ARG:NH2	3:A:602:HOH:O	2.28	0.57
1:C:344:ARG:NH2	1:C:349:ASP:OD1	2.38	0.57
1:D:340:LEU:HD13	1:D:343:GLN:OE1	2.05	0.57
1:D:165:MET:HE1	1:D:201:LEU:C	2.25	0.56
1:D:268:HIS:ND1	1:D:270:ILE:HG12	2.20	0.56
1:A:169:ALA:HB2	1:A:205:TYR:HB3	1.87	0.56
1:A:158:VAL:HG12	1:B:158:VAL:HG12	1.87	0.55
1:D:253:PHE:HE1	1:D:288:ILE:HD12	1.71	0.55
1:A:25:PHE:HE2	1:A:53:VAL:HG21	1.71	0.54
1:C:31:MET:HE3	1:C:37:MET:HG3	1.90	0.54
1:A:363:GLU:H	1:A:363:GLU:CD	2.11	0.54
1:A:17[B]:ARG:NH2	1:A:136:GLU:OE1	2.39	0.54
1:D:308:ILE:O	1:D:312:ARG:HG3	2.08	0.53
1:C:63:THR:HB	1:C:89:ILE:HB	1.90	0.53
1:C:412:PRO:HG3	1:D:5:ARG:HD3	1.89	0.53
1:C:252:ARG:NH2	3:C:601:HOH:O	2.00	0.52
1:D:380:ASP:OD1	1:D:382:LYS:HE3	2.09	0.52
1:A:146:VAL:HB	1:A:168:ILE:HG12	1.91	0.52
1:B:149:TYR:HD2	1:B:405:GLU:HG2	1.75	0.52
1:D:304:VAL:O	1:D:308:ILE:HG13	2.09	0.52
1:B:380:ASP:OD2	1:B:382:LYS:NZ	2.35	0.51
1:A:81:ARG:O	1:A:83:PRO:HD3	2.11	0.51
1:C:5:ARG:O	1:C:30:ARG:NH1	2.43	0.51
1:D:216:ILE:HB	1:D:258:ILE:HD12	1.93	0.51
1:C:379:LEU:HB3	1:C:425:LEU:HD13	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:298:PRO:HG2	1:C:301:HIS:HB2	1.93	0.51
1:C:262:SER:HB2	1:C:266:ILE:HB	1.92	0.50
1:D:358:VAL:O	1:D:362:ILE:N	2.44	0.50
1:B:260:VAL:HG11	1:B:265:LEU:HG	1.92	0.50
1:D:385:PRO:HB2	1:D:388:PRO:HG2	1.92	0.50
1:A:1:MET:HG3	1:A:162:GLU:HG2	1.93	0.50
1:D:9:GLN:NE2	1:D:30:ARG:HG2	2.26	0.50
1:D:42:ILE:HG22	1:D:44:GLY:H	1.75	0.50
1:C:149:TYR:OH	1:C:407:GLU:OE2	2.22	0.50
1:D:63:THR:HB	1:D:89:ILE:HB	1.93	0.49
1:D:352:HIS:O	1:D:352:HIS:ND1	2.46	0.49
1:D:299:ILE:HG23	1:D:361:VAL:HG11	1.95	0.49
1:B:344:ARG:HG3	1:B:345:PRO:O	2.14	0.48
1:D:31:MET:SD	1:D:32:PRO:HD2	2.53	0.48
1:D:354:ASP:OD1	1:D:356:ASN:ND2	2.46	0.48
1:A:45:ILE:HA	1:A:108:GLU:O	2.14	0.48
1:C:327:ILE:HB	1:C:353:VAL:HG21	1.95	0.48
1:D:165:MET:CE	1:D:201:LEU:HB3	2.43	0.48
1:C:18:GLU:HG3	1:C:23:THR:HG21	1.95	0.47
1:A:377:GLU:H	1:A:377:GLU:CD	2.18	0.47
1:B:368:ILE:HG13	1:B:398:TYR:CZ	2.50	0.47
1:B:241[B]:ARG:HH22	1:D:78:ASP:HB2	1.80	0.47
1:C:171:LEU:HD21	1:C:420:ARG:HA	1.97	0.47
1:D:148:LEU:HB2	1:D:183:HIS:CE1	2.49	0.47
1:A:63:THR:HB	1:A:89:ILE:HB	1.96	0.47
1:B:264:ASP:HB2	1:B:301:HIS:HE2	1.79	0.47
1:B:371:LYS:HE3	1:B:371:LYS:HB2	1.62	0.46
1:C:294:HIS:HD2	1:C:295:SER:N	2.12	0.46
1:A:220:LEU:HA	1:B:41:LEU:HD21	1.97	0.46
1:A:391:LYS:HD2	1:A:395:ASP:OD2	2.15	0.46
1:D:141:PRO:HB2	1:D:399:THR:HG22	1.96	0.46
1:A:329:GLN:HG3	1:A:383:GLN:HB3	1.98	0.46
1:D:311:ILE:HG12	1:D:319:PHE:HD2	1.81	0.46
1:D:344:ARG:HG3	1:D:348:LEU:HD12	1.96	0.46
1:D:354:ASP:HB3	1:D:357:ASP:OD1	2.16	0.46
1:A:165:MET:CE	1:A:202:LEU:HD23	2.46	0.46
1:C:377:GLU:OE1	1:C:377:GLU:N	2.40	0.46
1:A:304:VAL:HG12	1:A:308:ILE:HD12	1.98	0.45
1:C:357:ASP:OD2	3:C:602:HOH:O	2.21	0.45
1:D:300:LYS:HD3	1:D:300:LYS:HA	1.75	0.45
1:B:311:ILE:HG13	1:B:319:PHE:HD2	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:PRO:HB3	1:A:347:PHE:HB3	1.97	0.45
1:C:216:ILE:HD13	1:C:274:VAL:HG11	1.99	0.45
1:A:25:PHE:CE2	1:A:53:VAL:HG21	2.51	0.45
1:B:149:TYR:HE2	1:B:407:GLU:HG3	1.82	0.45
1:C:356:ASN:O	1:C:359:PHE:HD1	1.99	0.45
1:A:17[A]:ARG:HG2	1:A:22:ILE:HG12	1.99	0.45
1:B:1:MET:HA	1:B:163:THR:OG1	2.16	0.45
1:C:232:VAL:HG12	1:C:270:ILE:HB	1.98	0.44
1:D:15:ASN:HB2	1:D:136:GLU:OE2	2.18	0.44
1:D:253:PHE:HD1	1:D:288:ILE:HB	1.82	0.44
1:D:340:LEU:HB2	1:D:343:GLN:HG3	1.98	0.44
1:C:358:VAL:O	1:C:362:ILE:HB	2.18	0.44
1:A:159:MET:HE1	1:A:416:ILE:HD11	2.00	0.44
1:B:310:LEU:HD13	1:B:313:ARG:NH2	2.33	0.44
1:C:17[A]:ARG:NH1	1:C:20:GLY:O	2.51	0.44
1:C:177:GLU:HG3	1:C:211:ASN:O	2.18	0.44
1:C:181:GLU:HB3	1:D:81:ARG:HH22	1.82	0.44
1:D:152:MET:HE3	1:D:410:ARG:NH2	2.33	0.44
1:D:179:LEU:HD13	1:D:215:TRP:HH2	1.83	0.44
1:D:267:PRO:HB2	1:D:271:TRP:CD1	2.52	0.43
1:D:324:ASP:OD2	1:D:371:LYS:NZ	2.50	0.43
1:C:144:TYR:HE1	1:C:426:ILE:HG13	1.83	0.43
1:C:162:GLU:HG3	1:C:201:LEU:HD22	2.00	0.43
1:C:257:LYS:HA	1:C:257:LYS:HD3	1.86	0.43
1:C:63:THR:HA	1:C:67:GLN:O	2.18	0.43
1:D:181:GLU:N	3:D:618:HOH:O	2.50	0.43
1:A:262:SER:OG	1:A:266:ILE:HD12	2.18	0.43
1:A:17[B]:ARG:HH22	1:A:136:GLU:CD	2.19	0.43
1:A:143:ARG:O	1:A:401:TYR:HA	2.19	0.43
1:C:17[B]:ARG:NH2	1:C:136:GLU:OE1	2.52	0.43
1:C:301:HIS:CE1	1:C:303:VAL:HG13	2.54	0.43
1:B:81:ARG:O	1:B:83:PRO:HD3	2.18	0.43
1:A:216:ILE:HD11	1:A:239:LEU:HB2	2.00	0.43
1:B:63:THR:HB	1:B:89:ILE:HB	2.00	0.43
1:C:3:LEU:O	3:C:603:HOH:O	2.22	0.43
1:B:236:ALA:HA	1:B:274:VAL:HG13	2.01	0.42
1:B:268:HIS:ND1	1:B:270:ILE:HG12	2.34	0.42
1:B:282:ALA:HB1	1:B:287:VAL:O	2.18	0.42
1:A:282:ALA:HB1	1:A:287:VAL:O	2.18	0.42
1:C:15:ASN:HB2	1:C:136:GLU:OE2	2.19	0.42
1:C:355:PRO:HG2	1:C:388:PRO:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:358:VAL:HA	1:D:361:VAL:HG22	2.02	0.42
1:C:266:ILE:HD13	1:C:303:VAL:HG12	2.02	0.42
1:D:26:GLN:HA	1:D:90:ILE:O	2.19	0.42
1:A:126:ARG:HH11	1:A:126:ARG:HG3	1.85	0.42
1:B:241[B]:ARG:NH2	1:D:78:ASP:HB2	2.35	0.42
1:B:241[B]:ARG:NH1	1:D:78:ASP:OD2	2.46	0.42
1:C:344:ARG:HD2	1:C:348:LEU:HD12	2.01	0.42
1:B:139:GLY:HA3	1:B:427:ARG:HG2	2.01	0.42
1:B:362:ILE:HD12	1:B:362:ILE:HA	1.93	0.42
1:D:325:THR:OG1	1:D:371:LYS:N	2.52	0.42
1:B:409:GLU:OE2	1:B:411:GLU:HG2	2.19	0.42
1:C:5:ARG:NH2	1:D:412:PRO:HA	2.35	0.42
1:C:31:MET:SD	1:C:32:PRO:HD2	2.59	0.41
1:C:179:LEU:HB2	1:C:182:ALA:HB3	2.00	0.41
1:C:296:PRO:HB3	1:C:347:PHE:HB3	2.02	0.41
1:B:412:PRO:HB2	1:B:413:TRP:CE3	2.55	0.41
1:C:34:TYR:HB3	1:D:152:MET:HE2	2.01	0.41
1:A:165:MET:SD	1:A:201:LEU:HB3	2.61	0.41
1:B:165:MET:HE1	1:B:202:LEU:CA	2.46	0.41
1:A:1:MET:CG	1:A:162:GLU:HG2	2.50	0.41
1:A:82:TRP:CE2	1:A:88:ALA:HB2	2.55	0.41
1:A:149:TYR:OH	1:A:407:GLU:HG3	2.21	0.41
1:A:257:LYS:HB2	1:A:257:LYS:HE3	1.73	0.41
1:B:310:LEU:HD13	1:B:313:ARG:HH21	1.86	0.41
1:B:345:PRO:HD2	1:B:348:LEU:HD12	2.02	0.41
1:C:45:ILE:HA	1:C:108:GLU:O	2.21	0.41
1:C:34:TYR:CD2	1:D:149:TYR:HE2	2.39	0.40
1:D:240:GLN:O	1:D:244:ARG:HG3	2.20	0.40
1:A:182:ALA:O	1:B:84:LEU:HD23	2.21	0.40
1:D:174:THR:O	1:D:208:GLU:HG2	2.21	0.40
1:A:57:PRO:HG3	1:A:110:ARG:HH11	1.87	0.40
1:A:3:LEU:HD13	1:A:30:ARG:HB2	2.04	0.40
1:C:107:ILE:HD11	1:C:127:VAL:HG22	2.03	0.40
1:C:171:LEU:HD23	1:C:171:LEU:HA	1.77	0.40
1:D:34:TYR:HH	1:D:115:TYR:HH	1.66	0.40
1:D:103:HIS:HA	3:D:647:HOH:O	2.20	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	434/435 (100%)	411 (95%)	23 (5%)	0	100	100
1	B	434/435 (100%)	414 (95%)	20 (5%)	0	100	100
1	C	434/435 (100%)	413 (95%)	21 (5%)	0	100	100
1	D	433/435 (100%)	409 (94%)	24 (6%)	0	100	100
All	All	1735/1740 (100%)	1647 (95%)	88 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	372/371 (100%)	366 (98%)	6 (2%)	58	64
1	B	372/371 (100%)	365 (98%)	7 (2%)	52	57
1	C	372/371 (100%)	359 (96%)	13 (4%)	31	31
1	D	371/371 (100%)	359 (97%)	12 (3%)	34	35
All	All	1487/1484 (100%)	1449 (97%)	38 (3%)	41	44

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ARG
1	A	119	GLU

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Mol	Chain	Res	Type
1	A	131	VAL
1	A	228	ARG
1	A	229	ASP
1	A	257	LYS
1	B	43	ASP
1	B	229	ASP
1	B	252	ARG
1	B	257	LYS
1	B	261	VAL
1	B	371	LYS
1	B	411	GLU
1	C	5	ARG
1	C	34	TYR
1	C	43	ASP
1	C	78	ASP
1	C	120	HIS
1	C	131	VAL
1	C	149	TYR
1	C	208	GLU
1	C	262	SER
1	C	263	SER
1	C	317	LYS
1	C	336	LYS
1	C	371	LYS
1	D	1	MET
1	D	21	GLU
1	D	43	ASP
1	D	229	ASP
1	D	252	ARG
1	D	255	ARG
1	D	262	SER
1	D	300	LYS
1	D	335	LEU
1	D	371	LYS
1	D	382	LYS
1	D	409	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	352	HIS
1	C	294	HIS

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Mol	Chain	Res	Type
1	C	352	HIS
1	D	356	ASN

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	435/435 (100%)	0.22	6 (1%) 73 72	20, 31, 48, 85	1 (0%)
1	B	435/435 (100%)	0.32	26 (5%) 29 27	18, 32, 56, 95	1 (0%)
1	C	435/435 (100%)	0.37	11 (2%) 58 57	22, 33, 51, 80	1 (0%)
1	D	435/435 (100%)	0.62	27 (6%) 28 26	24, 37, 58, 78	0
All	All	1740/1740 (100%)	0.38	70 (4%) 43 41	18, 33, 55, 95	3 (0%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	341	PRO	4.7
1	B	345	PRO	4.7
1	B	337	PRO	4.5
1	B	340	LEU	4.3
1	A	340	LEU	4.1
1	D	341	PRO	4.0
1	C	265	LEU	4.0
1	B	264	ASP	3.9
1	D	265	LEU	3.9
1	D	337	PRO	3.7
1	D	347	PHE	3.6
1	C	340	LEU	3.4
1	B	346	ALA	3.1
1	A	337	PRO	3.1
1	A	120	HIS	3.0
1	D	429	ILE	3.0
1	C	341	PRO	2.9
1	B	20	GLY	2.9
1	B	342	GLY	2.9
1	D	271	TRP	2.9
1	C	337	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	20	GLY	2.9
1	D	361	VAL	2.8
1	B	334	PRO	2.8
1	B	294	HIS	2.8
1	D	400	GLY	2.8
1	B	265	LEU	2.7
1	B	261	VAL	2.7
1	B	266	ILE	2.7
1	C	264	ASP	2.7
1	B	336	LYS	2.6
1	B	335	LEU	2.6
1	D	262	SER	2.6
1	C	118	GLN	2.6
1	A	366	VAL	2.6
1	C	271	TRP	2.5
1	C	342	GLY	2.5
1	D	346	ALA	2.5
1	B	361	VAL	2.5
1	B	435	GLU	2.4
1	D	435	GLU	2.4
1	D	359	PHE	2.4
1	B	333	ILE	2.4
1	B	263	SER	2.4
1	B	347	PHE	2.4
1	D	340	LEU	2.4
1	D	356	ASN	2.4
1	D	430	ALA	2.3
1	B	262	SER	2.3
1	B	352	HIS	2.3
1	D	434	ALA	2.3
1	C	352	HIS	2.3
1	D	260	VAL	2.2
1	B	348	LEU	2.2
1	A	263	SER	2.2
1	B	225	PRO	2.2
1	B	338	GLY	2.2
1	C	303	VAL	2.2
1	B	339	GLU	2.1
1	C	338	GLY	2.1
1	D	19	GLY	2.1
1	D	303	VAL	2.1
1	D	394	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	264	ASP	2.1
1	D	354	ASP	2.1
1	A	352	HIS	2.1
1	D	263	SER	2.1
1	D	267	PRO	2.0
1	D	392	ALA	2.0
1	D	327	ILE	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 monosaccharides in this entry.

6.4 Ligands [i](#)

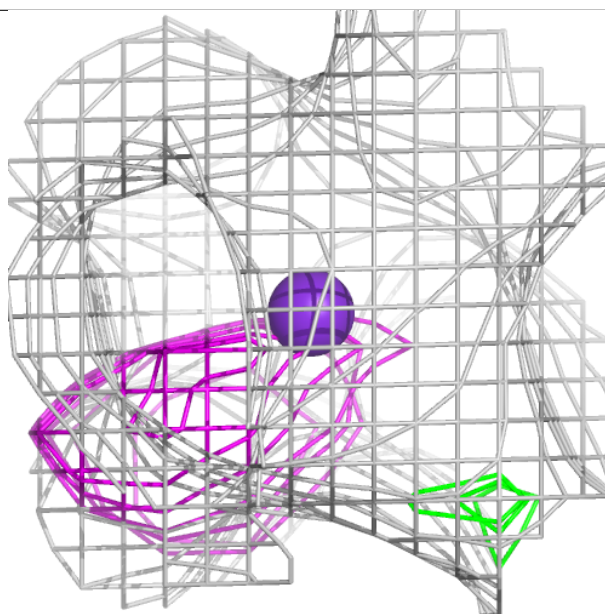
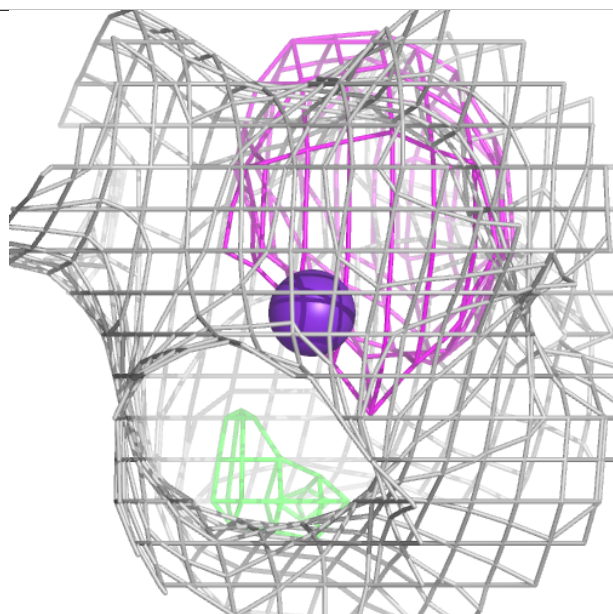
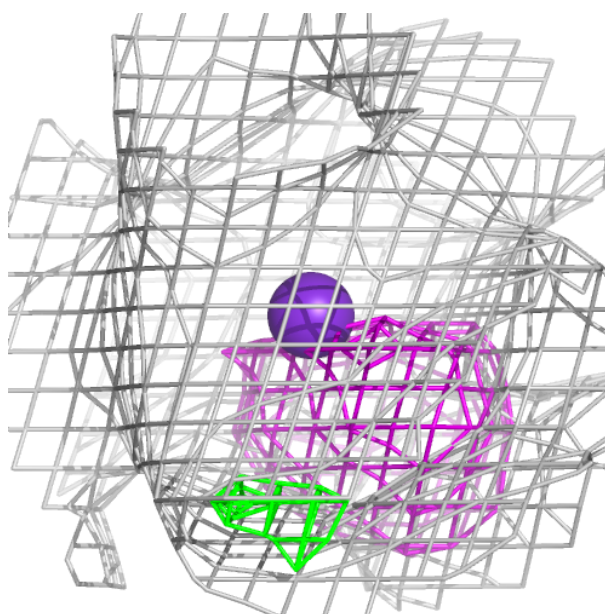
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	RB	B	501	1/1	0.97	0.09	52,52,52,52	0
2	RB	C	501	1/1	0.98	0.12	53,53,53,53	0
2	RB	D	501	1/1	0.98	0.14	66,66,66,66	0
2	RB	A	501	1/1	0.99	0.11	47,47,47,47	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.

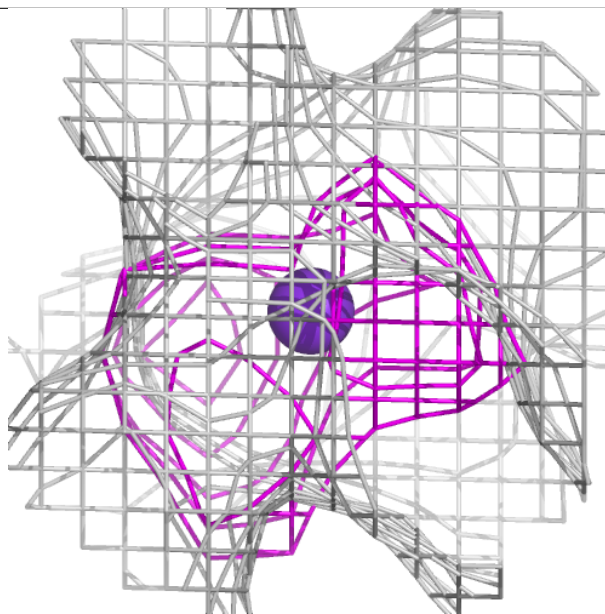
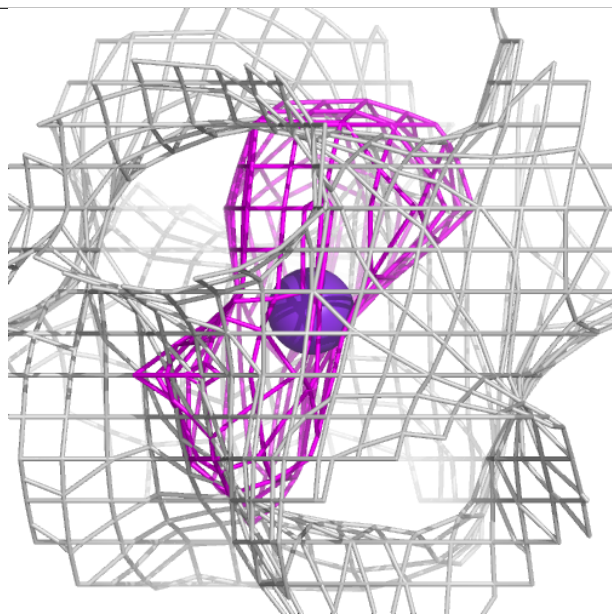
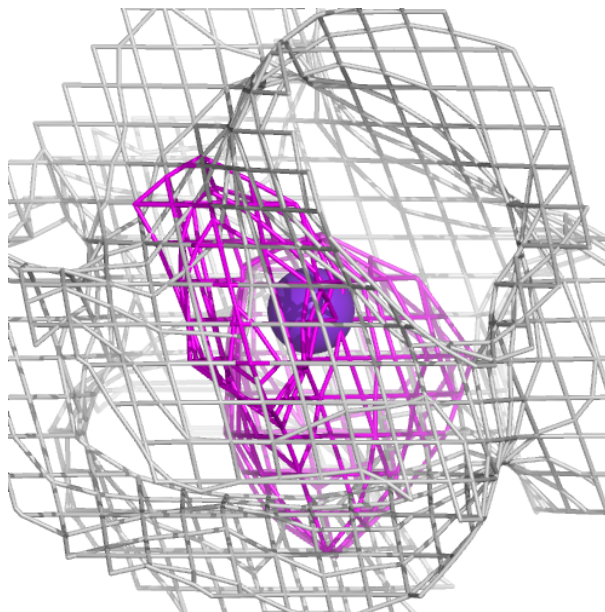
Electron density around RB B 501:

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



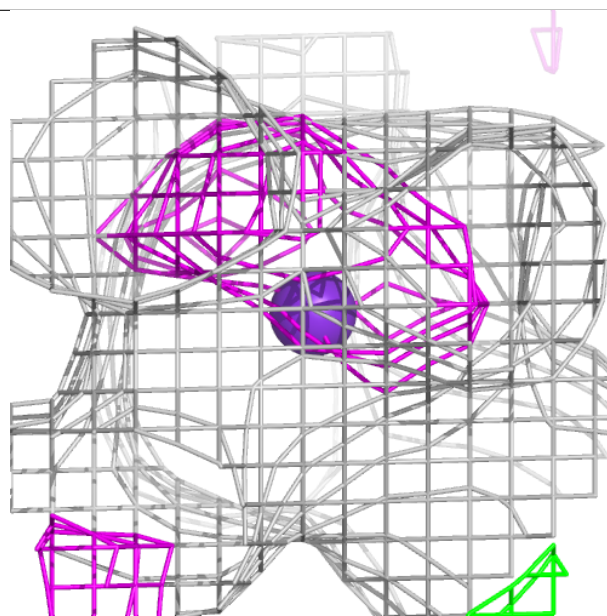
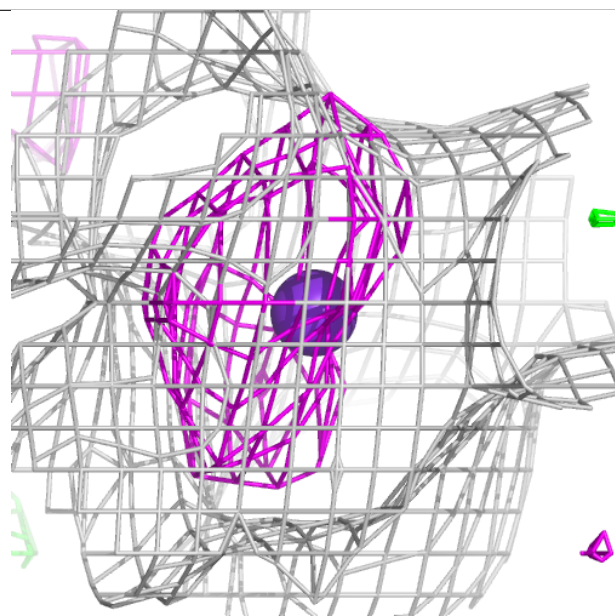
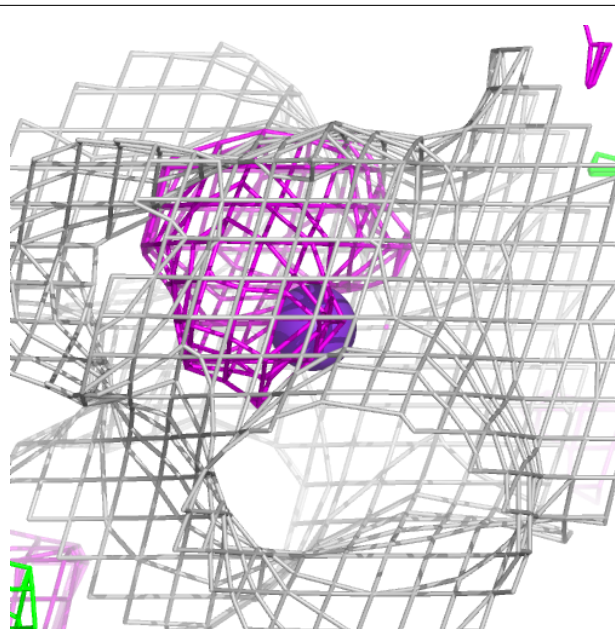
Electron density around RB C 501:

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



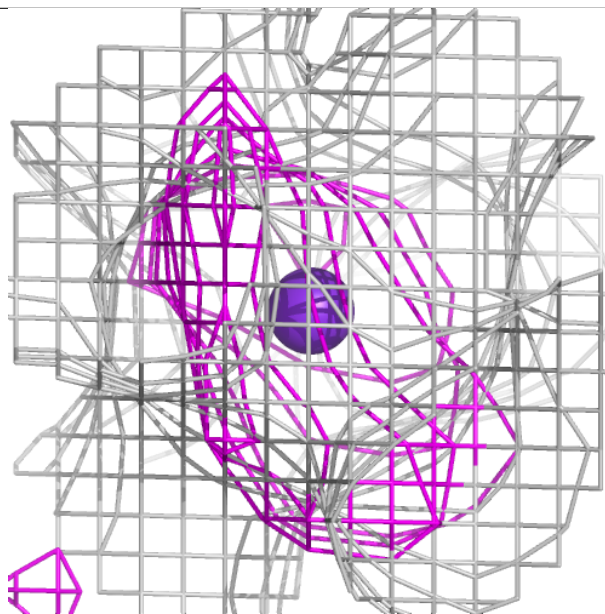
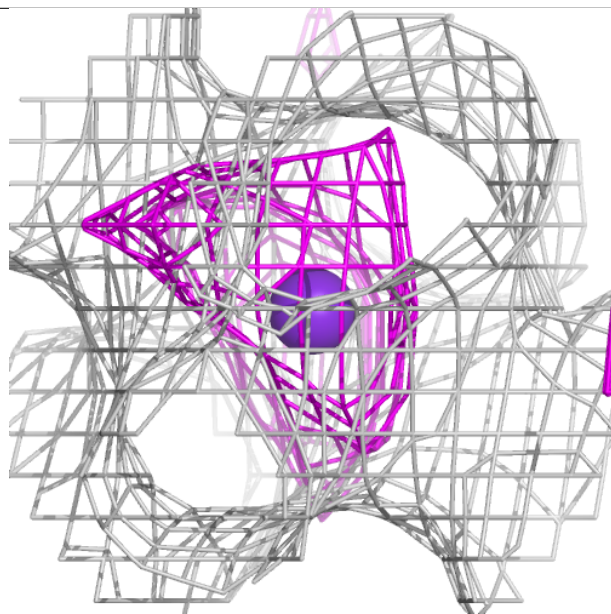
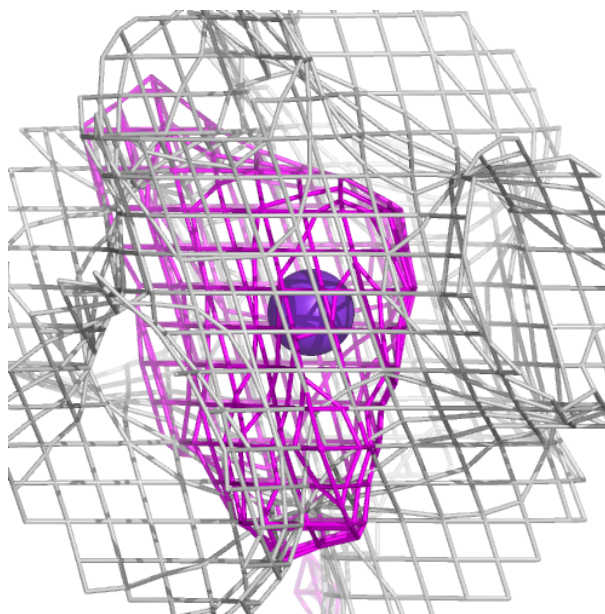
Electron density around RB D 501:

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



Electron density around RB A 501:

$2mF_o-DF_c$ (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.