



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 20, 2025 – 03:19 PM JST

PDB ID : 9K8U / pdb\_00009k8u  
Title : Ethanol dehydrogenase of Bursaphelenchus xylophilus  
Authors : Xing, Z.F.; Luo, X.; Li, X.Y.; Song, R.J.; Song, B.A.  
Deposited on : 2024-10-24  
Resolution : 2.81 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](mailto: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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
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)	:	2.46

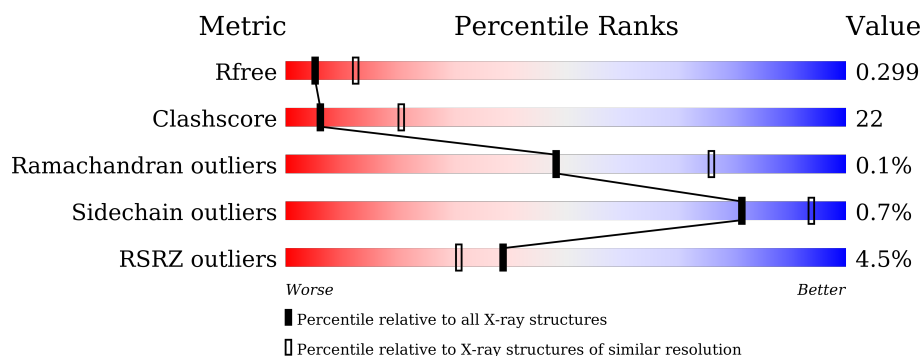
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.81 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	4293 (2.84-2.80)
Clashscore	180529	4801 (2.84-2.80)
Ramachandran outliers	177936	4739 (2.84-2.80)
Sidechain outliers	177891	4741 (2.84-2.80)
RSRZ outliers	164620	4295 (2.84-2.80)

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  $\geq 3$ , 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	347	<div> <div>5%</div> <div>67%</div> <div>31%</div> <div>.</div> </div>
1	B	347	<div> <div>4%</div> <div>61%</div> <div>38%</div> <div>.</div> </div>
1	C	347	<div> <div>5%</div> <div>59%</div> <div>39%</div> <div>.</div> </div>
1	D	347	<div> <div>5%</div> <div>57%</div> <div>42%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	B	402	-	-	X	-

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10504 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 alcohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	0	0
			2626	1664	454	491	17			
1	B	347	Total	C	N	O	S	0	0	0
			2627	1664	454	492	17			
1	C	347	Total	C	N	O	S	0	0	0
			2626	1664	454	491	17			
1	D	346	Total	C	N	O	S	0	0	0
			2617	1658	452	490	17			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

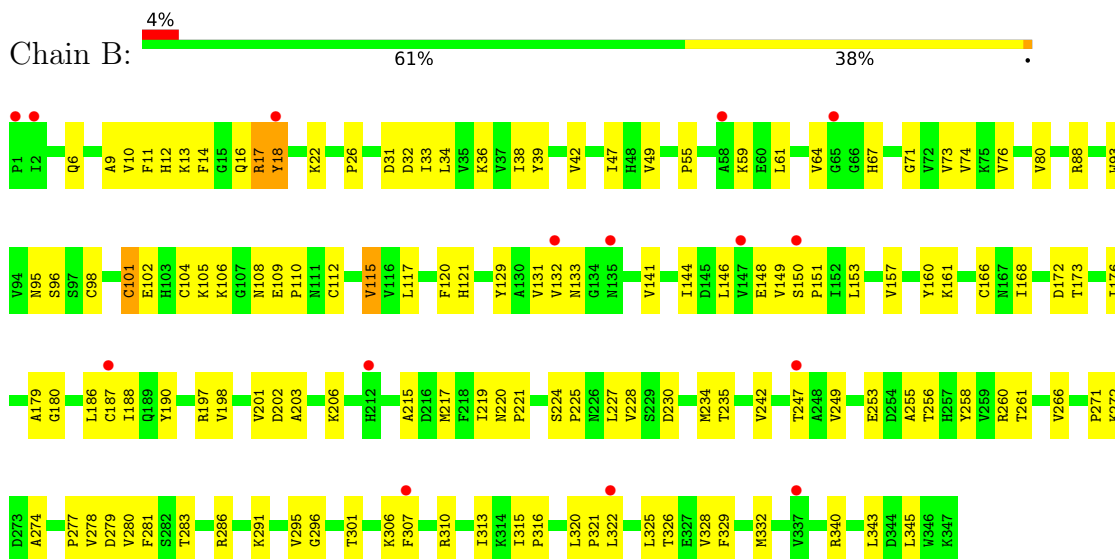
### 3 Residue-property plots [i](#)

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.

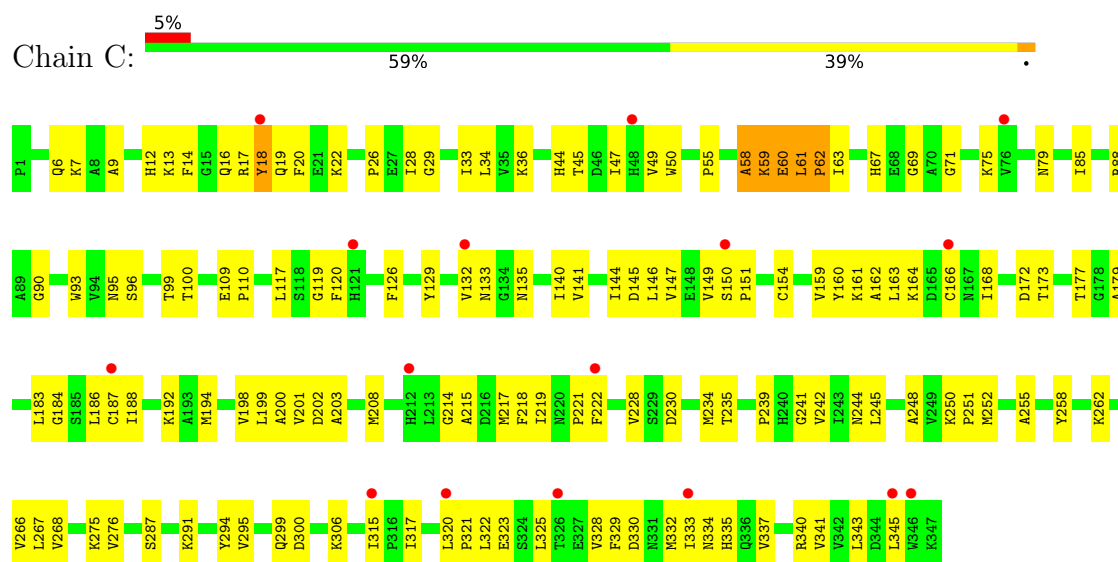
- Molecule 1: alcohol dehydrogenase



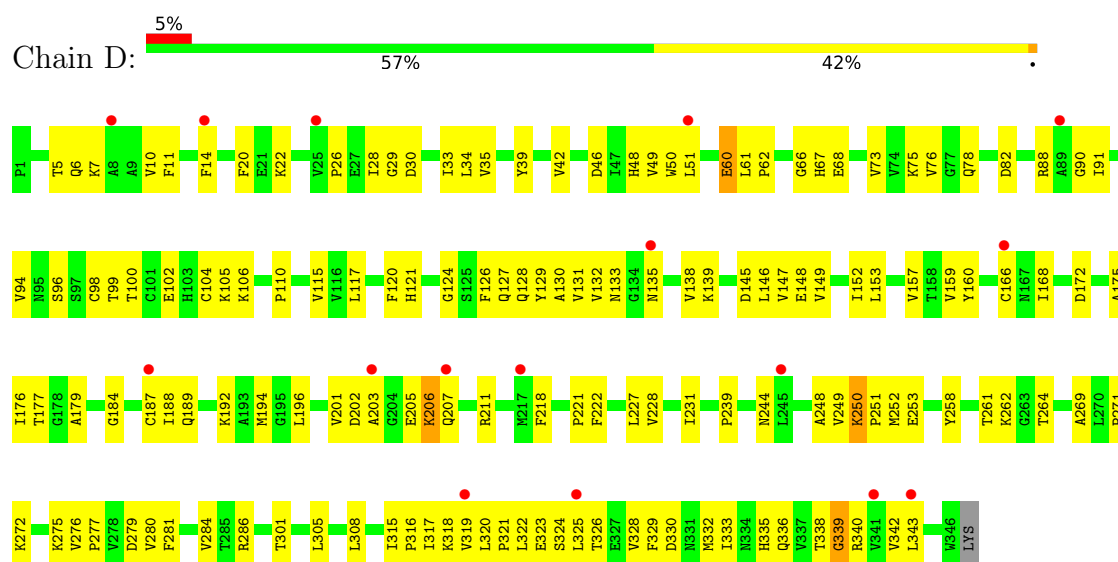
- Molecule 1: alcohol dehydrogenase



- Molecule 1: alcohol dehydrogenase



• Molecule 1: alcohol dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.71Å 152.77Å 63.79Å 90.00° 95.73° 90.00°	Depositor
Resolution (Å)	43.10 – 2.81 43.10 – 2.81	Depositor EDS
% Data completeness (in resolution range)	98.9 (43.10-2.81) 98.8 (43.10-2.81)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.13 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.258 , 0.298 0.258 , 0.299	Depositor DCC
$R_{free}$ test set	1488 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	75.4	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 73.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	10504	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN

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	0.28	1/2676 (0.0%)	0.53	2/3627 (0.1%)
1	B	0.25	0/2677	0.56	2/3627 (0.1%)
1	C	0.29	0/2676	0.68	7/3627 (0.2%)
1	D	0.34	0/2667	0.66	3/3616 (0.1%)
All	All	0.29	1/10696 (0.0%)	0.61	14/14497 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	62	PRO	N-CD	7.05	1.57	1.47

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	ARG	N-CA-C	-12.26	91.69	109.96
1	C	58	ALA	N-CA-C	-10.66	93.79	109.15
1	C	208	MET	N-CA-C	-10.11	100.94	113.28
1	C	62	PRO	N-CA-C	-10.05	94.73	110.50
1	C	59	LYS	N-CA-C	-8.86	102.05	113.12
1	D	206	LYS	N-CA-C	-8.49	102.92	113.20
1	B	17	ARG	CB-CA-C	8.19	121.95	110.16
1	A	61	LEU	C-N-CD	-8.04	92.02	125.00
1	C	60	GLU	N-CA-C	-7.15	99.30	109.96
1	A	211	ARG	CG-CD-NE	-6.88	96.87	112.00
1	C	61	LEU	N-CA-C	-6.42	95.62	109.81
1	D	339	GLY	N-CA-C	-5.71	106.05	112.33
1	D	60	GLU	N-CA-C	-5.37	105.83	112.38
1	C	61	LEU	CB-CA-C	5.01	120.03	110.17

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	2626	0	2647	89	0
1	B	2627	0	2651	127	0
1	C	2626	0	2647	124	0
1	D	2617	0	2637	147	0
2	A	2	0	0	0	0
2	B	2	0	0	2	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
All	All	10504	0	10582	464	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 22.

All (464) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ILE:HA	1:B:345:LEU:HD22	1.48	0.96
1:B:98:CYS:SG	2:B:402:ZN:ZN	1.57	0.94
1:A:10:VAL:HG22	1:A:64:VAL:HG22	1.53	0.90
1:A:91:ILE:HD12	1:A:132:VAL:HG21	1.55	0.88
1:B:104:CYS:SG	1:B:112:CYS:HB2	2.12	0.88
1:C:28:ILE:HD13	1:C:75:LYS:HD3	1.54	0.87
1:C:179:ALA:HB1	1:C:188:ILE:HD11	1.57	0.87
1:B:161:LYS:HD2	1:B:296:GLY:HA2	1.60	0.83
1:D:250:LYS:HG3	1:D:251:PRO:HD3	1.60	0.83
1:B:328:VAL:HG12	1:B:332:MET:HE3	1.59	0.83
1:D:11:PHE:O	1:D:62:PRO:HA	1.78	0.83
1:B:112:CYS:HB3	1:B:115:VAL:HG23	1.60	0.83
1:B:16:GLN:HG3	1:B:17:ARG:O	1.78	0.82
1:A:147:VAL:HG22	1:A:317:ILE:CD1	2.10	0.81
1:A:44:HIS:HA	1:A:332:MET:HE3	1.60	0.81
1:B:55:PRO:HD3	1:B:272:LYS:NZ	1.95	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ASN:HD21	1:D:30:ASP:HB2	1.46	0.80
1:B:179:ALA:HB1	1:B:188:ILE:HD11	1.62	0.80
1:B:148:GLU:HA	1:B:315:ILE:HD11	1.65	0.79
1:B:76:VAL:HG22	1:B:80:VAL:HG21	1.66	0.78
1:B:256:THR:HG21	1:B:278:VAL:CG2	2.14	0.78
1:C:252:MET:HE3	1:C:267:LEU:HD22	1.66	0.77
1:A:91:ILE:HG23	1:A:132:VAL:HG21	1.67	0.76
1:A:168:ILE:HD11	1:A:241:GLY:HA3	1.67	0.76
1:B:11:PHE:HB3	1:B:18:TYR:HB3	1.67	0.75
1:B:102:GLU:HA	1:B:105:LYS:HE2	1.67	0.74
1:D:176:ILE:CD1	1:D:187:CYS:SG	2.76	0.74
1:A:122:ARG:HE	1:A:131:VAL:HG23	1.52	0.74
1:A:70:ALA:HB1	1:A:146:LEU:HD22	1.69	0.74
1:D:28:ILE:HG22	1:D:29:GLY:O	1.88	0.73
1:D:325:LEU:CD2	1:D:343:LEU:HD11	2.17	0.73
1:D:325:LEU:HG	1:D:325:LEU:O	1.89	0.73
1:D:166:CYS:HB3	1:D:168:ILE:HG13	1.71	0.73
1:A:166:CYS:HB3	1:A:168:ILE:HG13	1.70	0.73
1:D:177:THR:HG22	1:D:244:ASN:HA	1.70	0.72
1:B:151:PRO:HG2	1:B:315:ILE:HD13	1.70	0.72
1:D:177:THR:CG2	1:D:244:ASN:HA	2.19	0.72
1:D:194:MET:HB2	1:D:196:LEU:HG	1.71	0.72
1:D:94:VAL:HG21	1:D:110:PRO:HG3	1.70	0.72
1:B:326:THR:HA	1:B:329:PHE:HD2	1.54	0.71
1:D:96:SER:HB3	1:D:133:ASN:HB2	1.71	0.71
1:C:99:THR:HG22	1:C:99:THR:O	1.89	0.70
1:D:325:LEU:HD22	1:D:343:LEU:CD2	2.22	0.70
1:C:79:ASN:HD21	1:D:30:ASP:CB	2.04	0.70
1:C:168:ILE:HD11	1:C:241:GLY:HA3	1.74	0.70
1:C:28:ILE:HD11	1:C:34:LEU:HD22	1.73	0.70
1:C:192:LYS:HD2	1:C:214:GLY:HA3	1.74	0.70
1:D:325:LEU:HD22	1:D:343:LEU:HD21	1.74	0.69
1:A:277:PRO:HB3	1:D:275:LYS:HD2	1.74	0.69
1:B:96:SER:HB3	1:B:133:ASN:HB2	1.75	0.68
1:A:12:HIS:ND1	1:A:62:PRO:HG3	2.08	0.68
1:A:147:VAL:HG22	1:A:317:ILE:HD11	1.76	0.68
1:D:7:LYS:HG2	1:D:22:LYS:HG3	1.75	0.68
1:B:256:THR:HG21	1:B:278:VAL:HG23	1.75	0.68
1:C:61:LEU:O	1:C:63:ILE:HG22	1.93	0.68
1:D:221:PRO:HA	1:D:227:LEU:HD21	1.75	0.68
1:A:47:ILE:HG13	1:A:332:MET:HG2	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:ILE:HD11	1:D:187:CYS:SG	2.34	0.68
1:A:179:ALA:HB1	1:A:188:ILE:HD11	1.74	0.67
1:B:101:CYS:HB2	1:B:104:CYS:SG	2.33	0.67
1:C:47:ILE:HG12	1:C:332:MET:HB2	1.75	0.67
1:B:307:PHE:HB3	1:B:313:ILE:HD13	1.76	0.67
1:A:43:CYS:HB3	1:A:68:GLU:OE2	1.95	0.67
1:D:20:PHE:CE1	1:D:323:GLU:HG2	2.30	0.67
1:A:91:ILE:HD12	1:A:132:VAL:CG2	2.23	0.67
1:A:152:ILE:HG13	1:A:308:LEU:HD22	1.77	0.66
1:D:261:THR:HG22	1:D:286:ARG:HG2	1.77	0.66
1:D:22:LYS:H	1:D:22:LYS:HD2	1.60	0.66
1:A:325:LEU:HD13	1:A:343:LEU:HD11	1.76	0.66
1:B:256:THR:CG2	1:B:278:VAL:HG23	2.26	0.65
1:A:211:ARG:HD3	1:A:215:ALA:O	1.95	0.65
1:B:88:ARG:HD3	1:B:146:LEU:HD11	1.78	0.65
1:D:325:LEU:CD2	1:D:343:LEU:CD1	2.75	0.65
1:A:147:VAL:HG22	1:A:317:ILE:HD12	1.77	0.65
1:D:5:THR:HB	1:D:22:LYS:HB3	1.78	0.64
1:B:277:PRO:HB2	1:C:275:LYS:HD3	1.79	0.64
1:C:60:GLU:HG2	1:C:62:PRO:O	1.97	0.64
1:C:228:VAL:HG13	1:C:258:TYR:HB3	1.80	0.64
1:D:91:ILE:HG12	1:D:132:VAL:HG11	1.77	0.64
1:B:166:CYS:HB3	1:B:168:ILE:HG12	1.78	0.64
1:D:179:ALA:HB1	1:D:188:ILE:HD11	1.78	0.64
1:A:21:GLU:HG2	1:A:22:LYS:N	2.11	0.64
1:A:308:LEU:HD12	1:A:313:ILE:O	1.97	0.64
1:C:28:ILE:HG12	1:C:34:LEU:HB2	1.79	0.63
1:B:6:GLN:O	1:B:22:LYS:HA	1.97	0.63
1:B:47:ILE:CD1	1:B:329:PHE:HD1	2.12	0.63
1:D:61:LEU:HD22	1:D:61:LEU:H	1.62	0.63
1:B:325:LEU:HD22	1:B:343:LEU:HD11	1.81	0.63
1:C:187:CYS:SG	1:C:245:LEU:HD11	2.39	0.62
1:C:100:THR:HB	1:D:82:ASP:OD2	1.99	0.62
1:C:325:LEU:HD13	1:C:343:LEU:HD11	1.82	0.62
1:A:242:VAL:HG11	1:A:255:ALA:HB1	1.80	0.62
1:B:180:GLY:HA3	1:B:206:LYS:HD3	1.81	0.62
1:C:88:ARG:HD2	1:C:146:LEU:HD11	1.82	0.61
1:A:315:ILE:HD12	1:A:316:PRO:HD2	1.82	0.61
1:D:88:ARG:HD3	1:D:146:LEU:HD21	1.82	0.61
1:A:47:ILE:HG13	1:A:332:MET:CG	2.29	0.61
1:B:117:LEU:H	1:B:121:HIS:HB2	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:VAL:HG13	1:D:276:VAL:O	2.01	0.61
1:D:319:VAL:HA	1:D:342:VAL:HG13	1.83	0.61
1:D:332:MET:HA	1:D:336:GLN:H	1.66	0.61
1:C:337:VAL:HG21	1:C:340:ARG:HH12	1.65	0.60
1:B:326:THR:HA	1:B:329:PHE:CD2	2.36	0.60
1:D:320:LEU:O	1:D:343:LEU:HA	2.02	0.60
1:D:33:ILE:HG22	1:D:132:VAL:O	2.02	0.59
1:B:34:LEU:HD11	1:B:129:TYR:HB3	1.84	0.59
1:C:28:ILE:CG1	1:C:34:LEU:HB2	2.31	0.59
1:A:183:LEU:HD13	1:A:245:LEU:HD21	1.82	0.59
1:C:126:PHE:O	1:C:322:LEU:HD11	2.03	0.59
1:A:168:ILE:HG23	1:A:172:ASP:HB2	1.85	0.59
1:C:50:TRP:HZ3	1:C:333:ILE:HD12	1.66	0.59
1:D:250:LYS:CG	1:D:251:PRO:HD3	2.33	0.59
1:B:98:CYS:SG	1:B:101:CYS:HB2	2.42	0.59
1:B:256:THR:CG2	1:B:278:VAL:CG2	2.79	0.58
1:C:163:LEU:HB3	1:C:194:MET:HE1	1.83	0.58
1:A:176:ILE:HD12	1:A:187:CYS:HB3	1.85	0.58
1:B:217:MET:HG2	1:B:234:MET:HE2	1.86	0.58
1:C:299:GLN:HA	1:D:105:LYS:HD2	1.86	0.58
1:C:14:PHE:CZ	1:C:58:ALA:HB3	2.39	0.58
1:B:230:ASP:O	1:B:234:MET:HG2	2.03	0.58
1:A:308:LEU:CD1	1:A:313:ILE:O	2.52	0.57
1:D:175:ALA:HB2	1:D:239:PRO:HG2	1.86	0.57
1:A:91:ILE:HG23	1:A:132:VAL:CG2	2.32	0.57
1:D:35:VAL:HG22	1:D:73:VAL:HG22	1.86	0.57
1:D:316:PRO:C	1:D:317:ILE:HG13	2.27	0.57
1:A:10:VAL:HG22	1:A:64:VAL:CG2	2.31	0.57
1:D:321:PRO:HD2	1:D:324:SER:HB2	1.86	0.57
1:B:93:TRP:CD1	1:B:295:VAL:HG21	2.40	0.57
1:D:248:ALA:O	1:D:251:PRO:HD2	2.05	0.57
1:C:166:CYS:HB3	1:C:168:ILE:HG13	1.87	0.56
1:C:79:ASN:ND2	1:D:30:ASP:HB2	2.19	0.56
1:C:332:MET:HE2	1:C:337:VAL:HG12	1.88	0.56
1:C:135:ASN:HA	1:D:100:THR:CG2	2.35	0.56
1:D:159:VAL:HG11	1:D:187:CYS:HA	1.87	0.56
1:C:13:LYS:HB2	1:C:16:GLN:HB2	1.87	0.56
1:C:96:SER:HB3	1:C:133:ASN:HB2	1.88	0.56
1:B:328:VAL:O	1:B:332:MET:HG3	2.06	0.56
1:C:334:ASN:O	1:C:335:HIS:CG	2.59	0.56
1:D:152:ILE:O	1:D:157:VAL:HG23	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:242:VAL:HG11	1:C:255:ALA:HB1	1.88	0.55
1:C:183:LEU:HB3	1:C:245:LEU:HD22	1.88	0.55
1:B:261:THR:HG23	1:B:286:ARG:HA	1.88	0.55
1:D:98:CYS:O	1:D:99:THR:OG1	2.24	0.55
1:A:9:ALA:HA	1:A:19:GLN:O	2.06	0.55
1:B:11:PHE:HZ	1:B:49:VAL:HG12	1.72	0.55
1:B:98:CYS:HG	2:B:402:ZN:ZN	1.15	0.55
1:C:322:LEU:HB2	1:C:345:LEU:HA	1.88	0.55
1:D:176:ILE:HD13	1:D:187:CYS:HB3	1.88	0.55
1:D:325:LEU:HD22	1:D:343:LEU:CD1	2.36	0.54
1:A:103:HIS:HA	1:A:106:LYS:HE2	1.89	0.54
1:D:20:PHE:CZ	1:D:323:GLU:HG2	2.42	0.54
1:D:33:ILE:HG13	1:D:76:VAL:HG12	1.90	0.54
1:A:255:ALA:HA	1:A:258:TYR:CZ	2.42	0.54
1:D:10:VAL:HG11	1:D:62:PRO:HB3	1.90	0.54
1:D:168:ILE:HG23	1:D:172:ASP:HB2	1.90	0.54
1:D:127:GLN:NE2	1:D:130:ALA:HA	2.22	0.54
1:D:179:ALA:HA	1:D:184:GLY:HA3	1.90	0.54
1:A:317:ILE:HA	1:A:340:ARG:O	2.08	0.54
1:C:49:VAL:HG11	1:C:119:GLY:O	2.08	0.54
1:B:12:HIS:CD2	1:B:12:HIS:H	2.27	0.53
1:B:328:VAL:CG1	1:B:332:MET:HE3	2.35	0.53
1:B:55:PRO:HD3	1:B:272:LYS:HZ1	1.72	0.53
1:C:28:ILE:HD12	1:C:28:ILE:H	1.73	0.53
1:B:188:ILE:HG23	1:B:198:VAL:HG11	1.89	0.53
1:D:317:ILE:HA	1:D:340:ARG:HB2	1.89	0.53
1:A:177:THR:O	1:A:245:LEU:HB3	2.09	0.53
1:C:198:VAL:HB	1:C:215:ALA:HA	1.90	0.53
1:B:272:LYS:HE2	1:B:272:LYS:HA	1.91	0.53
1:B:316:PRO:HD2	1:B:340:ARG:HH21	1.74	0.52
1:C:330:ASP:O	1:C:333:ILE:HG22	2.10	0.52
1:C:276:VAL:HG13	1:C:276:VAL:O	2.08	0.52
1:B:39:TYR:HA	1:B:343:LEU:O	2.10	0.52
1:C:99:THR:O	1:C:99:THR:CG2	2.58	0.52
1:B:18:TYR:HH	1:B:329:PHE:HZ	1.55	0.52
1:C:203:ALA:HB1	1:C:222:PHE:CE1	2.45	0.52
1:D:46:ASP:HA	1:D:49:VAL:HG23	1.91	0.52
1:A:18:TYR:HH	1:A:50:TRP:CG	2.28	0.52
1:B:202:ASP:O	1:B:221:PRO:HD2	2.08	0.52
1:C:12:HIS:O	1:C:61:LEU:HG	2.10	0.52
1:D:326:THR:HA	1:D:329:PHE:HB3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:PRO:HD3	1:B:272:LYS:HZ3	1.73	0.52
1:B:101:CYS:H	1:B:104:CYS:HB2	1.75	0.52
1:C:159:VAL:HB	1:C:187:CYS:HB3	1.92	0.52
1:A:206:LYS:HB2	1:A:218:PHE:CZ	2.45	0.52
1:A:276:VAL:HG23	1:A:276:VAL:O	2.10	0.52
1:B:98:CYS:SG	1:B:112:CYS:SG	3.07	0.52
1:D:176:ILE:HD13	1:D:187:CYS:SG	2.49	0.51
1:D:339:GLY:O	1:D:340:ARG:C	2.49	0.51
1:B:280:VAL:CG2	1:C:252:MET:HE1	2.41	0.51
1:D:67:HIS:CE1	1:D:68:GLU:HG2	2.46	0.51
1:A:39:TYR:HA	1:A:343:LEU:O	2.10	0.51
1:D:319:VAL:HG12	1:D:342:VAL:CG1	2.40	0.51
1:A:108:ASN:HB2	1:D:262:LYS:NZ	2.25	0.51
1:B:17:ARG:NH1	1:B:17:ARG:HB2	2.26	0.51
1:B:320:LEU:O	1:B:343:LEU:HA	2.10	0.51
1:C:45:THR:O	1:C:49:VAL:HG23	2.10	0.51
1:D:110:PRO:O	1:D:115:VAL:HG11	2.11	0.51
1:A:203:ALA:HB2	1:A:221:PRO:HG2	1.93	0.51
1:A:138:VAL:CG1	1:A:305:LEU:HD11	2.41	0.51
1:B:74:VAL:HG21	1:B:129:TYR:CE2	2.45	0.51
1:C:135:ASN:HA	1:D:100:THR:HG21	1.92	0.51
1:D:328:VAL:O	1:D:332:MET:HG3	2.11	0.51
1:A:79:ASN:HD22	1:B:31:ASP:HB3	1.76	0.51
1:D:250:LYS:HG3	1:D:251:PRO:CD	2.37	0.51
1:B:10:VAL:HB	1:B:64:VAL:CG1	2.41	0.50
1:C:173:THR:HG21	1:C:235:THR:HG22	1.93	0.50
1:D:253:GLU:CD	1:D:277:PRO:HD2	2.36	0.50
1:A:109:GLU:N	1:A:110:PRO:HD2	2.26	0.50
1:D:202:ASP:OD1	1:D:206:LYS:HD2	2.12	0.50
1:C:334:ASN:O	1:C:335:HIS:ND1	2.45	0.50
1:A:210:CYS:HB3	1:A:215:ALA:HB3	1.94	0.50
1:D:39:TYR:HB3	1:D:342:VAL:CG2	2.42	0.50
1:D:317:ILE:HG23	1:D:340:ARG:HB3	1.93	0.50
1:A:185:SER:HB2	1:A:213:LEU:HD11	1.92	0.50
1:B:221:PRO:HA	1:B:227:LEU:HD21	1.93	0.50
1:D:317:ILE:HG23	1:D:340:ARG:CB	2.42	0.50
1:C:100:THR:HG23	1:D:139:LYS:HE2	1.93	0.50
1:D:42:VAL:HG21	1:D:325:LEU:HD21	1.93	0.50
1:B:9:ALA:HB1	1:B:18:TYR:CZ	2.47	0.50
1:C:266:VAL:HA	1:C:291:LYS:O	2.12	0.50
1:A:28:ILE:HD11	1:A:75:LYS:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:LYS:HB3	1:C:20:PHE:CE1	2.47	0.50
1:A:106:LYS:HE3	1:D:262:LYS:HD2	1.94	0.49
1:B:280:VAL:HG22	1:C:252:MET:HE1	1.95	0.49
1:C:67:HIS:CD2	1:C:154:CYS:SG	3.05	0.49
1:D:33:ILE:HD11	1:D:73:VAL:HG21	1.94	0.49
1:D:50:TRP:CD1	1:D:51:LEU:H	2.29	0.49
1:B:67:HIS:HB2	1:B:93:TRP:HA	1.94	0.49
1:C:255:ALA:HA	1:C:258:TYR:CZ	2.47	0.49
1:A:209:HIS:O	1:A:212:HIS:HB2	2.12	0.49
1:B:36:LYS:O	1:B:71:GLY:HA3	2.13	0.49
1:B:228:VAL:HG13	1:B:258:TYR:HB3	1.94	0.49
1:C:334:ASN:C	1:C:335:HIS:CG	2.91	0.49
1:A:45:THR:O	1:A:49:VAL:HG23	2.13	0.49
1:D:127:GLN:HE21	1:D:130:ALA:CB	2.24	0.49
1:A:138:VAL:HG13	1:A:305:LEU:HD11	1.93	0.49
1:A:252:MET:HB2	1:A:276:VAL:HG12	1.93	0.49
1:C:109:GLU:N	1:C:110:PRO:HD2	2.27	0.49
1:B:9:ALA:HB1	1:B:18:TYR:CE1	2.47	0.49
1:A:6:GLN:NE2	1:A:26:PRO:HD3	2.28	0.49
1:A:242:VAL:HB	1:A:265:VAL:HG12	1.93	0.49
1:D:189:GLN:O	1:D:192:LYS:HG2	2.13	0.49
1:B:325:LEU:HD22	1:B:343:LEU:CD1	2.43	0.48
1:A:253:GLU:OE2	1:A:276:VAL:HA	2.14	0.48
1:C:88:ARG:HB2	1:C:140:ILE:O	2.13	0.48
1:C:321:PRO:HB2	1:C:323:GLU:HG2	1.95	0.48
1:A:47:ILE:HG13	1:A:332:MET:HB3	1.95	0.48
1:D:75:LYS:HG2	1:D:76:VAL:N	2.29	0.48
1:B:141:VAL:HB	1:B:144:ILE:HD13	1.94	0.48
1:C:262:LYS:HD3	1:C:287:SER:HB2	1.95	0.48
1:D:33:ILE:HD11	1:D:73:VAL:HG11	1.96	0.48
1:D:152:ILE:HG12	1:D:308:LEU:HB2	1.95	0.48
1:D:342:VAL:O	1:D:343:LEU:HD23	2.14	0.48
1:A:49:VAL:HG22	1:A:120:PHE:CD2	2.49	0.48
1:C:33:ILE:HG12	1:C:132:VAL:O	2.13	0.48
1:A:12:HIS:HA	1:A:62:PRO:HA	1.96	0.48
1:A:49:VAL:HG22	1:A:120:PHE:HD2	1.78	0.48
1:A:319:VAL:HA	1:A:342:VAL:O	2.13	0.48
1:D:7:LYS:HB3	1:D:20:PHE:CZ	2.48	0.48
1:A:49:VAL:HG11	1:A:119:GLY:O	2.13	0.48
1:B:33:ILE:HD11	1:B:73:VAL:HG21	1.95	0.48
1:B:49:VAL:HG22	1:B:120:PHE:HD1	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ILE:HG23	1:B:172:ASP:HB2	1.95	0.48
1:C:49:VAL:HG22	1:C:120:PHE:HD1	1.78	0.48
1:B:47:ILE:HD11	1:B:329:PHE:HD1	1.77	0.47
1:B:74:VAL:HG21	1:B:129:TYR:HE2	1.79	0.47
1:D:269:ALA:C	1:D:271:PRO:HD3	2.39	0.47
1:B:203:ALA:HA	1:B:220:ASN:OD1	2.13	0.47
1:C:36:LYS:O	1:C:71:GLY:HA3	2.14	0.47
1:B:173:THR:HG23	1:B:197:ARG:HE	1.79	0.47
1:A:82:ASP:OD2	1:B:98:CYS:HB2	2.14	0.47
1:B:109:GLU:N	1:B:110:PRO:HD2	2.30	0.47
1:C:6:GLN:O	1:C:22:LYS:HA	2.13	0.47
1:D:6:GLN:HG3	1:D:6:GLN:O	2.15	0.47
1:A:33:ILE:HG23	1:A:73:VAL:HG13	1.96	0.47
1:C:18:TYR:CZ	1:C:329:PHE:HB3	2.50	0.47
1:C:49:VAL:HG22	1:C:120:PHE:CD1	2.49	0.47
1:A:112:CYS:HB3	1:A:115:VAL:HG13	1.96	0.47
1:B:279:ASP:O	1:B:283:THR:HG22	2.15	0.47
1:C:184:GLY:HA2	1:C:245:LEU:HD13	1.96	0.47
1:C:322:LEU:HD22	1:C:345:LEU:HG	1.97	0.47
1:C:150:SER:OG	1:C:151:PRO:HD3	2.15	0.47
1:A:252:MET:HE3	1:A:267:LEU:HD22	1.96	0.47
1:C:100:THR:CG2	1:D:139:LYS:HE2	2.45	0.47
1:C:14:PHE:CE1	1:C:58:ALA:HB3	2.50	0.46
1:D:7:LYS:HD2	1:D:128:GLN:NE2	2.30	0.46
1:C:145:ASP:O	1:C:149:VAL:HG23	2.15	0.46
1:C:100:THR:O	1:C:100:THR:HG22	2.14	0.46
1:B:108:ASN:HD22	1:C:262:LYS:HE3	1.81	0.46
1:C:47:ILE:HD13	1:C:332:MET:SD	2.56	0.46
1:C:147:VAL:HG22	1:C:317:ILE:HD12	1.98	0.46
1:D:33:ILE:HD11	1:D:73:VAL:CG1	2.46	0.46
1:D:330:ASP:O	1:D:333:ILE:HG22	2.16	0.46
1:C:117:LEU:HB2	1:C:120:PHE:HB3	1.98	0.46
1:C:184:GLY:HA2	1:C:245:LEU:CD1	2.45	0.46
1:C:230:ASP:O	1:C:234:MET:HG2	2.15	0.46
1:B:247:THR:HB	1:B:271:PRO:HB3	1.98	0.45
1:B:18:TYR:HD1	1:B:18:TYR:O	1.99	0.45
1:B:42:VAL:HG21	1:B:325:LEU:HD11	1.97	0.45
1:B:253:GLU:OE2	1:B:277:PRO:HD2	2.16	0.45
1:D:176:ILE:HD13	1:D:187:CYS:CB	2.45	0.45
1:C:160:TYR:OH	1:C:164:LYS:HE3	2.16	0.45
1:B:10:VAL:HB	1:B:64:VAL:HG12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:TRP:NE1	1:B:117:LEU:HD22	2.32	0.45
1:A:228:VAL:HA	1:A:258:TYR:HB3	1.99	0.45
1:A:256:THR:HG21	1:A:278:VAL:HB	1.99	0.45
1:B:6:GLN:HE22	1:B:26:PRO:HD3	1.81	0.45
1:B:146:LEU:HA	1:B:149:VAL:HG22	1.97	0.45
1:B:173:THR:HG21	1:B:235:THR:HG22	1.97	0.45
1:D:14:PHE:CE1	1:D:60:GLU:HB3	2.52	0.45
1:D:33:ILE:HD11	1:D:73:VAL:CG2	2.46	0.45
1:B:10:VAL:O	1:B:18:TYR:HB2	2.17	0.45
1:B:13:LYS:HE3	1:B:14:PHE:O	2.17	0.45
1:C:17:ARG:HA	1:C:17:ARG:HD3	1.61	0.45
1:B:160:TYR:CE1	1:B:307:PHE:HE2	2.35	0.45
1:B:281:PHE:CZ	1:C:55:PRO:HD2	2.52	0.45
1:D:20:PHE:O	1:D:20:PHE:CG	2.70	0.45
1:D:34:LEU:HD11	1:D:129:TYR:HB3	1.98	0.45
1:D:148:GLU:HA	1:D:317:ILE:CD1	2.46	0.45
1:B:321:PRO:O	1:B:343:LEU:HD22	2.17	0.44
1:B:322:LEU:HA	1:B:343:LEU:HD13	1.99	0.44
1:A:275:LYS:HG2	1:D:279:ASP:HA	1.99	0.44
1:B:176:ILE:HD13	1:B:187:CYS:HB3	1.99	0.44
1:D:318:LYS:N	1:D:340:ARG:O	2.50	0.44
1:A:229:SER:O	1:A:233:MET:HG2	2.16	0.44
1:B:249:VAL:O	1:B:253:GLU:HG2	2.17	0.44
1:C:199:LEU:HA	1:C:217:MET:O	2.18	0.44
1:D:166:CYS:HA	1:D:264:THR:HG21	1.99	0.44
1:B:266:VAL:HA	1:B:291:LYS:O	2.18	0.44
1:C:162:ALA:HB2	1:C:268:VAL:HG22	1.99	0.44
1:D:50:TRP:O	1:D:51:LEU:HB2	2.17	0.44
1:D:203:ALA:HB1	1:D:222:PHE:CZ	2.52	0.44
1:D:7:LYS:HB3	1:D:20:PHE:CE1	2.53	0.44
1:D:248:ALA:C	1:D:251:PRO:HD2	2.42	0.44
1:A:304:ALA:HA	1:A:307:PHE:HD2	1.83	0.44
1:B:255:ALA:HA	1:B:258:TYR:CZ	2.52	0.44
1:C:59:LYS:HA	1:C:59:LYS:HD3	1.73	0.44
1:D:90:GLY:N	1:D:138:VAL:HB	2.33	0.44
1:B:101:CYS:HB2	1:B:104:CYS:HG	1.83	0.44
1:C:325:LEU:HD22	1:C:343:LEU:HD11	2.00	0.44
1:D:48:HIS:NE2	1:D:272:LYS:HD3	2.33	0.44
1:C:151:PRO:HB2	1:C:315:ILE:HD12	2.00	0.44
1:A:47:ILE:HG13	1:A:332:MET:CB	2.48	0.43
1:D:6:GLN:CD	1:D:26:PRO:HD3	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:7:LYS:HD3	1:D:20:PHE:CZ	2.53	0.43
1:D:90:GLY:H	1:D:138:VAL:HB	1.82	0.43
1:C:28:ILE:HD12	1:C:28:ILE:N	2.33	0.43
1:B:59:LYS:HE2	1:B:61:LEU:HD22	1.99	0.43
1:B:280:VAL:HA	1:B:283:THR:HG22	2.00	0.43
1:C:252:MET:CE	1:C:267:LEU:HD22	2.42	0.43
1:A:68:GLU:HG2	1:A:154:CYS:HB2	2.00	0.43
1:D:147:VAL:O	1:D:317:ILE:HD13	2.19	0.43
1:A:265:VAL:CG2	1:A:290:VAL:HG22	2.48	0.43
1:C:93:TRP:CZ2	1:C:117:LEU:HD13	2.54	0.43
1:C:200:ALA:HB3	1:C:218:PHE:HD1	1.84	0.43
1:C:201:VAL:HG13	1:C:219:ILE:HG23	2.01	0.43
1:C:248:ALA:C	1:C:251:PRO:HD2	2.43	0.43
1:C:252:MET:HE2	1:C:252:MET:HB3	1.69	0.43
1:A:17:ARG:HG2	1:A:18:TYR:O	2.17	0.43
1:B:32:ASP:HB3	1:B:131:VAL:HG12	2.01	0.43
1:B:186:LEU:HD23	1:B:186:LEU:HA	1.72	0.43
1:C:161:LYS:HD3	1:C:300:ASP:OD2	2.17	0.43
1:C:250:LYS:HB3	1:C:250:LYS:HE3	1.77	0.43
1:D:66:GLY:O	1:D:124:GLY:HA3	2.18	0.43
1:C:6:GLN:NE2	1:C:26:PRO:HD3	2.34	0.43
1:B:190:TYR:CE2	1:B:313:ILE:HG12	2.54	0.43
1:C:328:VAL:HG21	1:C:341:VAL:HG11	2.00	0.43
1:D:145:ASP:O	1:D:149:VAL:HG23	2.18	0.43
1:B:13:LYS:HA	1:B:13:LYS:HD2	1.88	0.43
1:B:198:VAL:HB	1:B:215:ALA:HA	2.00	0.43
1:C:12:HIS:HD2	1:C:16:GLN:HE22	1.66	0.43
1:C:239:PRO:HD2	1:C:258:TYR:O	2.19	0.43
1:D:33:ILE:O	1:D:131:VAL:HA	2.18	0.43
1:C:306:LYS:HE2	1:C:306:LYS:HB3	1.74	0.43
1:D:253:GLU:OE1	1:D:277:PRO:HD2	2.19	0.43
1:A:11:PHE:CD2	1:A:50:TRP:HD1	2.37	0.42
1:A:157:VAL:HG22	1:A:301:THR:HA	2.01	0.42
1:A:238:GLY:C	1:A:260:ARG:HB2	2.44	0.42
1:A:267:LEU:HD13	1:D:280:VAL:HG13	2.01	0.42
1:B:17:ARG:HB2	1:B:17:ARG:CZ	2.48	0.42
1:C:18:TYR:HE2	1:C:50:TRP:CG	2.37	0.42
1:D:99:THR:HA	1:D:104:CYS:SG	2.59	0.42
1:C:177:THR:OG1	1:C:244:ASN:HA	2.19	0.42
1:D:102:GLU:O	1:D:106:LYS:HG3	2.18	0.42
1:D:332:MET:C	1:D:335:HIS:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:VAL:O	1:C:144:ILE:HG22	2.19	0.42
1:D:7:LYS:HE2	1:D:22:LYS:HE2	2.01	0.42
1:D:301:THR:O	1:D:305:LEU:HD23	2.20	0.42
1:D:319:VAL:HG12	1:D:342:VAL:HG13	2.00	0.42
1:D:325:LEU:HD22	1:D:343:LEU:HD11	1.94	0.42
1:C:50:TRP:CZ3	1:C:333:ILE:HD12	2.51	0.42
1:C:75:LYS:HA	1:C:85:ILE:HD11	2.01	0.42
1:D:6:GLN:OE1	1:D:26:PRO:HD3	2.18	0.42
1:A:228:VAL:O	1:A:232:GLN:HG3	2.20	0.42
1:A:252:MET:CB	1:A:276:VAL:HG12	2.49	0.42
1:B:157:VAL:HG22	1:B:301:THR:HA	2.01	0.42
1:B:201:VAL:HG22	1:B:219:ILE:CG2	2.49	0.42
1:C:14:PHE:HZ	1:C:58:ALA:HB3	1.81	0.42
1:A:14:PHE:CE1	1:A:58:ALA:HB3	2.55	0.42
1:B:148:GLU:O	1:B:151:PRO:HD2	2.20	0.42
1:D:30:ASP:HA	1:D:78:GLN:H	1.84	0.42
1:B:49:VAL:HG22	1:B:120:PHE:CD1	2.53	0.42
1:C:34:LEU:HD11	1:C:129:TYR:HB3	2.00	0.42
1:B:33:ILE:CD1	1:B:73:VAL:HG21	2.50	0.42
1:B:36:LYS:HB2	1:B:129:TYR:CE2	2.55	0.42
1:C:186:LEU:HD11	1:C:315:ILE:HG12	2.00	0.42
1:D:22:LYS:HD2	1:D:22:LYS:N	2.29	0.42
1:D:127:GLN:HE21	1:D:130:ALA:HA	1.83	0.42
1:D:159:VAL:CG1	1:D:187:CYS:HA	2.50	0.42
1:B:224:SER:HA	1:B:225:PRO:HD3	1.94	0.42
1:C:9:ALA:HA	1:C:19:GLN:O	2.19	0.42
1:A:185:SER:O	1:A:189:GLN:HG3	2.20	0.41
1:B:98:CYS:HB3	1:B:112:CYS:SG	2.59	0.41
1:A:331:ASN:HB3	1:A:336:GLN:HB2	2.02	0.41
1:B:98:CYS:SG	1:B:101:CYS:CB	3.09	0.41
1:B:242:VAL:HG11	1:B:255:ALA:HB1	2.01	0.41
1:C:47:ILE:HG13	1:C:329:PHE:HD1	1.85	0.41
1:C:201:VAL:HG22	1:C:219:ILE:CG2	2.50	0.41
1:D:117:LEU:HB2	1:D:120:PHE:HB2	2.02	0.41
1:D:121:HIS:O	1:D:121:HIS:CG	2.73	0.41
1:D:201:VAL:HG21	1:D:231:ILE:HD11	2.02	0.41
1:C:44:HIS:CD2	1:C:332:MET:HB3	2.56	0.41
1:C:202:ASP:O	1:C:221:PRO:HD2	2.21	0.41
1:D:6:GLN:NE2	1:D:26:PRO:HD3	2.35	0.41
1:B:249:VAL:HG22	1:B:274:ALA:HB1	2.01	0.41
1:C:28:ILE:HG22	1:C:29:GLY:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:TYR:HE1	1:D:194:MET:HE1	1.85	0.41
1:B:190:TYR:CZ	1:B:313:ILE:HG12	2.56	0.41
1:C:168:ILE:HG23	1:C:172:ASP:HB2	2.01	0.41
1:C:248:ALA:O	1:C:251:PRO:HD2	2.20	0.41
1:C:320:LEU:O	1:C:343:LEU:HA	2.20	0.41
1:D:126:PHE:HB3	1:D:322:LEU:HD11	2.02	0.41
1:D:205:GLU:C	1:D:207:GLN:H	2.29	0.41
1:D:207:GLN:HB2	1:D:218:PHE:CE1	2.55	0.41
1:A:94:VAL:HG22	1:A:110:PRO:HB3	2.02	0.41
1:A:269:ALA:O	1:D:284:VAL:HG21	2.21	0.41
1:D:207:GLN:HG3	1:D:211:ARG:HG2	2.03	0.41
1:A:94:VAL:CG2	1:A:110:PRO:HB3	2.50	0.41
1:B:106:LYS:HE3	1:B:106:LYS:HB2	1.83	0.41
1:B:150:SER:HA	1:B:153:LEU:HD12	2.02	0.41
1:B:151:PRO:HB3	1:B:340:ARG:HD3	2.03	0.41
1:C:99:THR:HG21	1:D:135:ASN:ND2	2.35	0.41
1:C:203:ALA:HB1	1:C:222:PHE:HE1	1.84	0.41
1:D:138:VAL:HG21	1:D:153:LEU:HG	2.03	0.41
1:D:152:ILE:HD12	1:D:315:ILE:HD12	2.03	0.41
1:D:228:VAL:HG13	1:D:258:TYR:HB3	2.02	0.41
1:D:336:GLN:O	1:D:338:THR:HG23	2.20	0.41
1:D:61:LEU:HB3	1:D:62:PRO:HD2	2.03	0.41
1:A:101:CYS:H	1:A:104:CYS:HB2	1.86	0.40
1:B:59:LYS:HG2	1:B:61:LEU:HB2	2.03	0.40
1:C:69:GLY:O	1:C:90:GLY:HA2	2.21	0.40
1:C:95:ASN:ND2	1:C:132:VAL:HG12	2.35	0.40
1:D:61:LEU:HD22	1:D:61:LEU:N	2.34	0.40
1:D:91:ILE:CG1	1:D:132:VAL:HG11	2.47	0.40
1:A:61:LEU:HG	1:A:61:LEU:O	2.21	0.40
1:B:306:LYS:HE3	1:B:310:ARG:HD3	2.03	0.40
1:D:279:ASP:OD2	1:D:281:PHE:HB3	2.21	0.40
1:C:294:TYR:CD2	1:C:295:VAL:HG22	2.55	0.40
1:D:207:GLN:HB2	1:D:218:PHE:CD1	2.56	0.40
1:B:33:ILE:HG22	1:B:132:VAL:O	2.21	0.40
1:B:47:ILE:HG13	1:B:332:MET:CB	2.51	0.40
1:D:249:VAL:O	1:D:253:GLU:HG2	2.21	0.40
1:B:95:ASN:ND2	1:B:132:VAL:HG22	2.36	0.40
1:B:260:ARG:HG2	1:B:261:THR:N	2.36	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	345/347 (99%)	332 (96%)	12 (4%)	1 (0%)	37	65
1	B	345/347 (99%)	336 (97%)	9 (3%)	0	100	100
1	C	345/347 (99%)	337 (98%)	8 (2%)	0	100	100
1	D	344/347 (99%)	327 (95%)	17 (5%)	0	100	100
All	All	1379/1388 (99%)	1332 (97%)	46 (3%)	1 (0%)	48	76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	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	289/289 (100%)	287 (99%)	2 (1%)	81	94
1	B	289/289 (100%)	286 (99%)	3 (1%)	73	91
1	C	289/289 (100%)	288 (100%)	1 (0%)	91	97
1	D	288/289 (100%)	286 (99%)	2 (1%)	81	94
All	All	1155/1156 (100%)	1147 (99%)	8 (1%)	81	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	CYS
1	A	68	GLU
1	B	18	TYR
1	B	101	CYS
1	B	115	VAL
1	C	18	TYR
1	D	250	LYS
1	D	252	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	111	ASN
1	A	207	GLN
1	B	103	HIS
1	B	108	ASN
1	C	12	HIS
1	C	79	ASN
1	C	95	ASN
1	C	103	HIS
1	D	6	GLN
1	D	121	HIS
1	D	127	GLN
1	D	128	GLN
1	D	135	ASN
1	D	207	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

Of 8 ligands modelled in this entry, 8 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

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	347/347 (100%)	0.39	16 (4%)	38 30	46, 73, 120, 148	0
1	B	347/347 (100%)	0.47	15 (4%)	40 33	43, 78, 146, 181	0
1	C	347/347 (100%)	0.35	16 (4%)	38 30	38, 67, 95, 134	0
1	D	346/347 (99%)	0.56	16 (4%)	38 30	46, 93, 176, 242	0
All	All	1387/1388 (99%)	0.44	63 (4%)	39 31	38, 75, 148, 242	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	166	CYS	3.8
1	A	108	ASN	3.7
1	A	166	CYS	3.6
1	A	328	VAL	3.5
1	A	343	LEU	3.3
1	B	18	TYR	3.2
1	C	166	CYS	3.1
1	B	322	LEU	3.1
1	C	132	VAL	3.0
1	C	187	CYS	3.0
1	A	207	GLN	3.0
1	C	18	TYR	3.0
1	A	247	THR	3.0
1	C	222	PHE	3.0
1	B	2	ILE	2.9
1	B	247	THR	2.9
1	B	212	HIS	2.8
1	A	312	GLN	2.8
1	B	187	CYS	2.8
1	B	135	ASN	2.8
1	C	346	TRP	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	203	ALA	2.7
1	D	14	PHE	2.6
1	C	48	HIS	2.6
1	D	217	MET	2.6
1	D	25	VAL	2.6
1	A	62	PRO	2.6
1	A	242	VAL	2.6
1	D	135	ASN	2.5
1	A	76	VAL	2.5
1	B	150	SER	2.5
1	A	58	ALA	2.5
1	D	207	GLN	2.4
1	C	121	HIS	2.4
1	D	89	ALA	2.4
1	D	245	LEU	2.4
1	D	319	VAL	2.4
1	B	307	PHE	2.4
1	D	325	LEU	2.3
1	A	269	ALA	2.3
1	D	187	CYS	2.3
1	C	320	LEU	2.3
1	C	345	LEU	2.2
1	B	147	VAL	2.2
1	B	337	VAL	2.2
1	C	333	ILE	2.2
1	C	150	SER	2.2
1	D	341	VAL	2.2
1	D	343	LEU	2.1
1	A	221	PRO	2.1
1	B	1	PRO	2.1
1	A	82	ASP	2.1
1	A	173	THR	2.1
1	C	315	ILE	2.1
1	B	132	VAL	2.1
1	D	8	ALA	2.1
1	A	28	ILE	2.1
1	B	65	GLY	2.1
1	C	212	HIS	2.1
1	C	326	THR	2.0
1	B	58	ALA	2.0
1	D	51	LEU	2.0
1	C	76	VAL	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 6.4 Ligands ⓘ

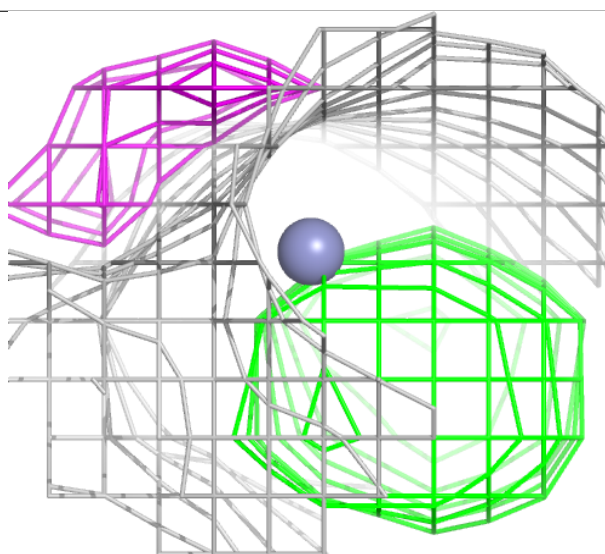
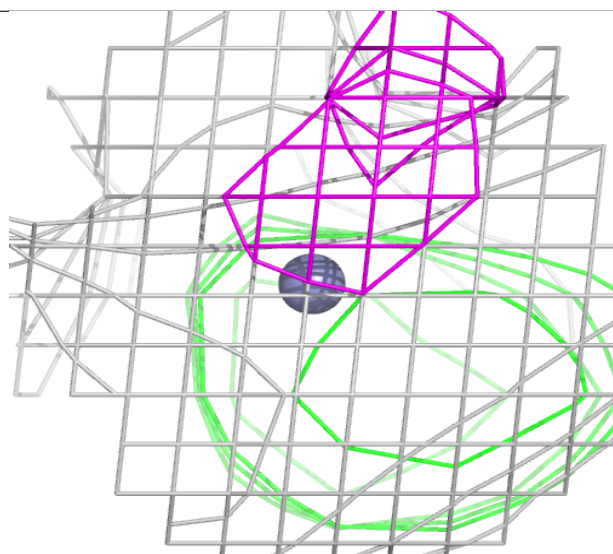
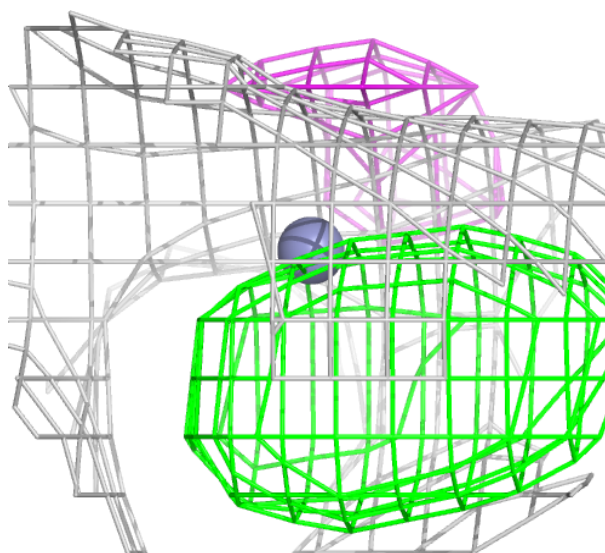
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	C	401	1/1	0.64	0.22	189,189,189,189	0
2	ZN	B	402	1/1	0.81	0.13	118,118,118,118	0
2	ZN	D	402	1/1	0.81	0.14	90,90,90,90	0
2	ZN	C	402	1/1	0.88	0.10	104,104,104,104	0
2	ZN	D	401	1/1	0.90	0.06	129,129,129,129	0
2	ZN	A	402	1/1	0.91	0.10	99,99,99,99	0
2	ZN	B	401	1/1	0.96	0.04	82,82,82,82	0
2	ZN	A	401	1/1	0.98	0.05	101,101,101,101	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.

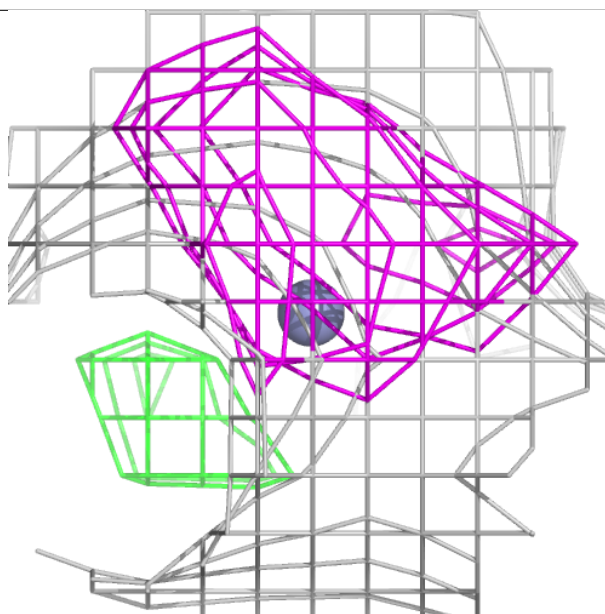
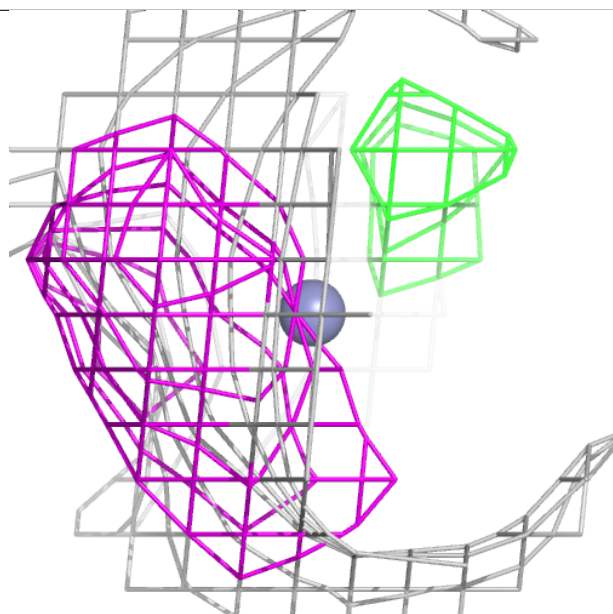
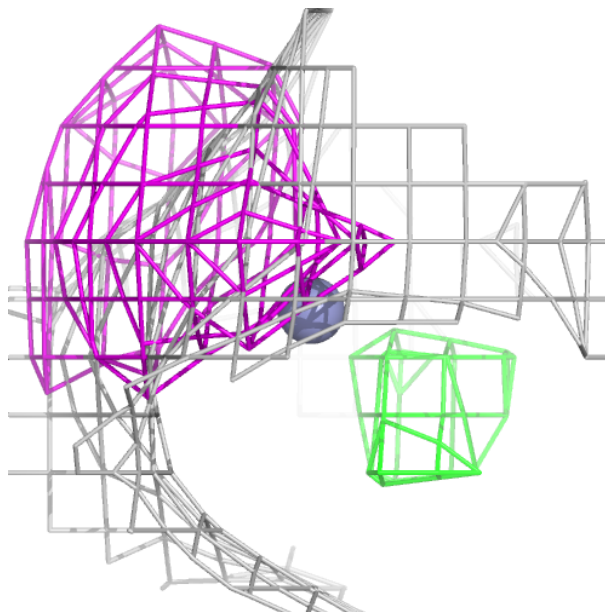
**Electron density around ZN C 401:**

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



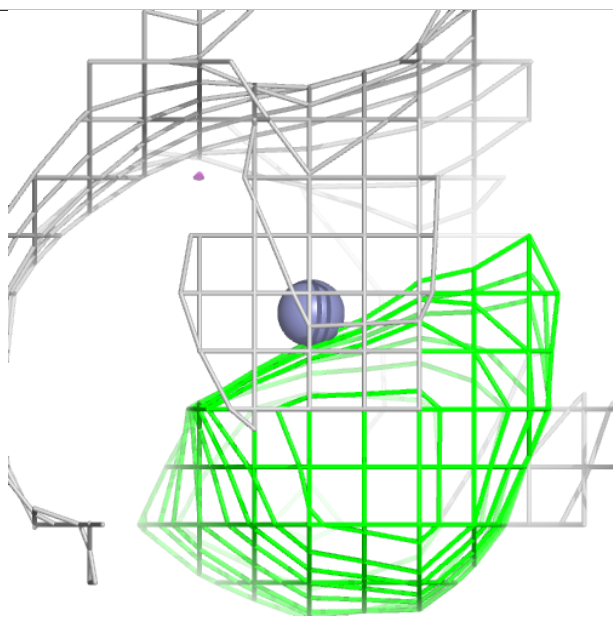
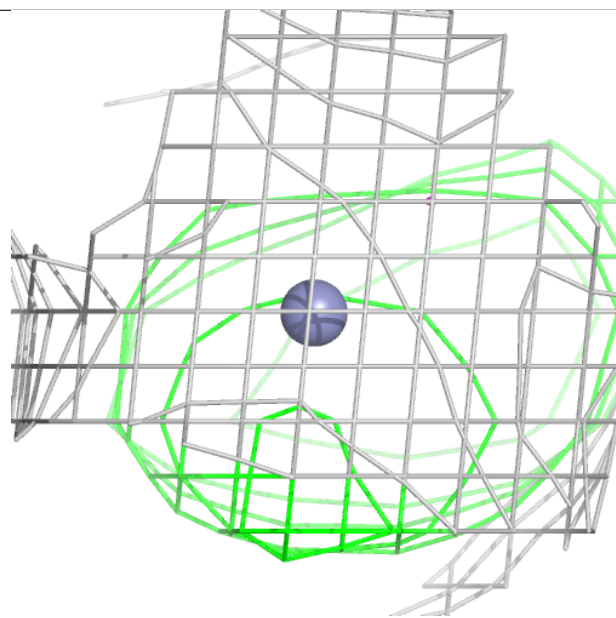
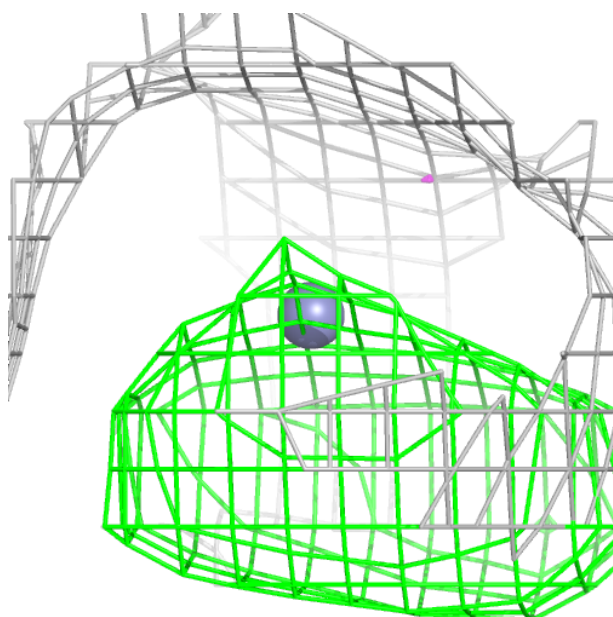
**Electron density around ZN B 402:**

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



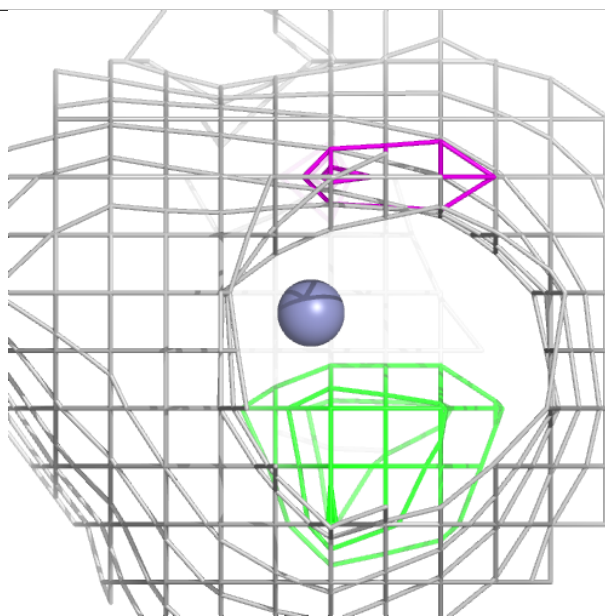
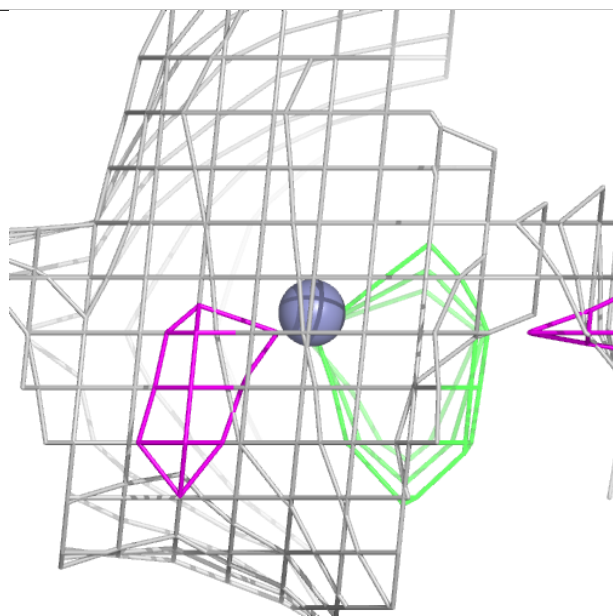
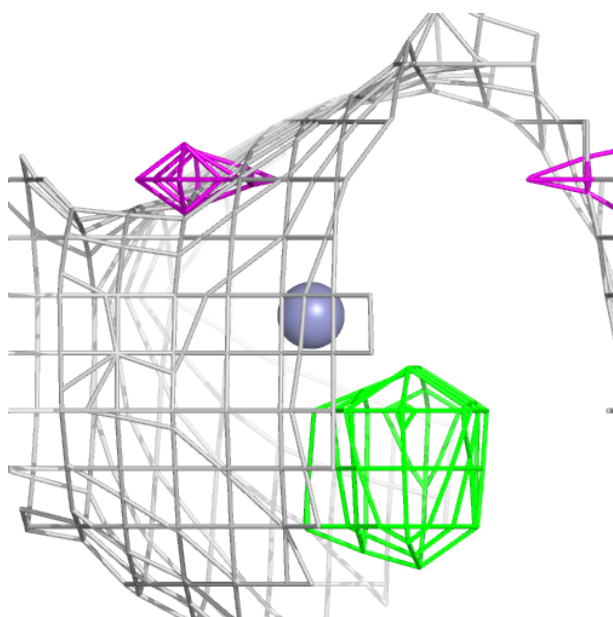
**Electron density around ZN D 402:**

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



**Electron density around ZN C 402:**

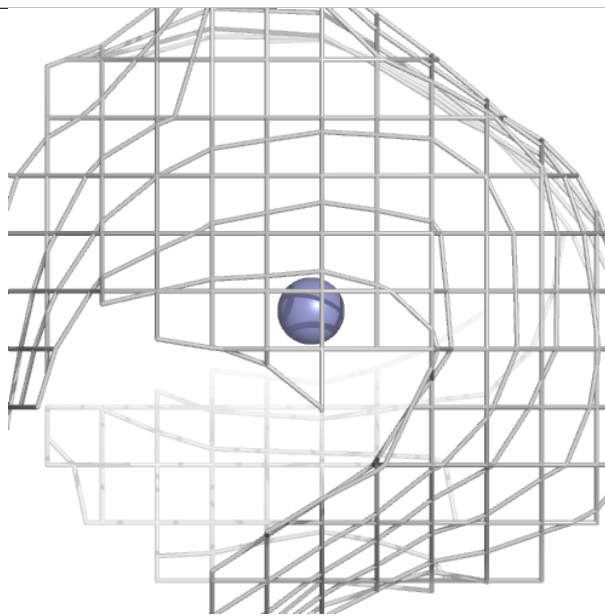
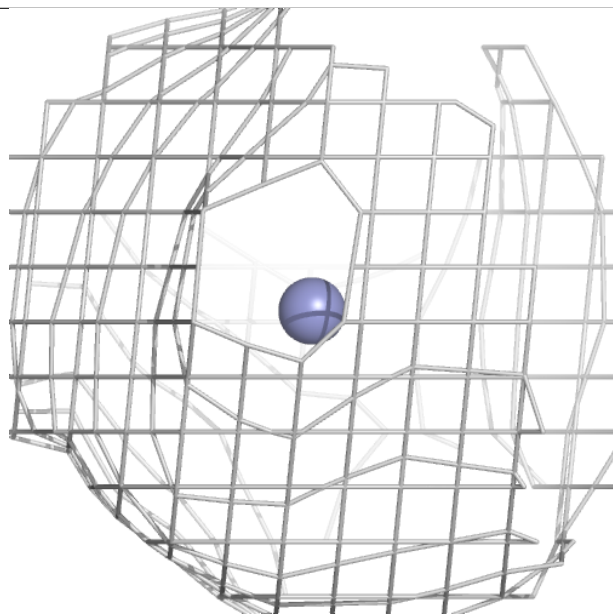
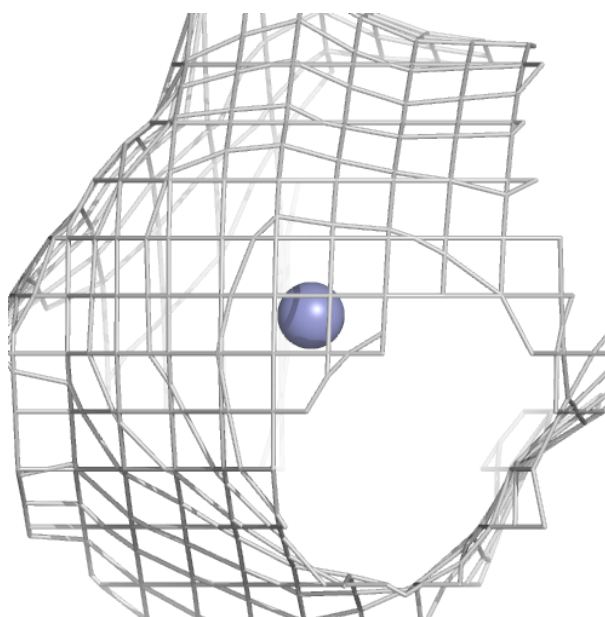
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





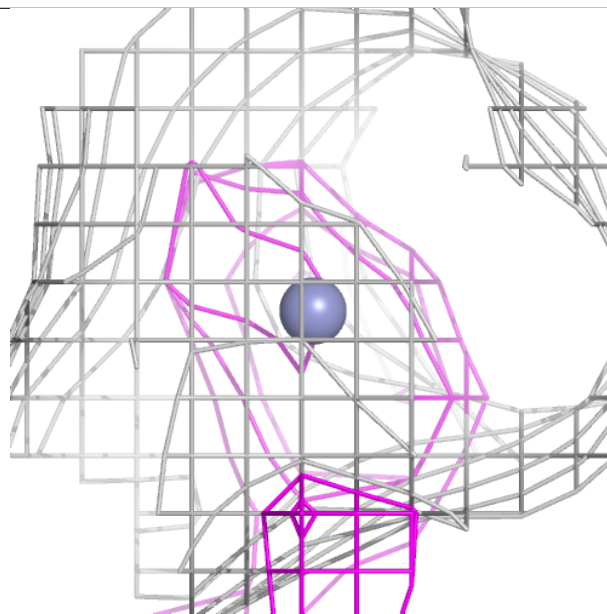
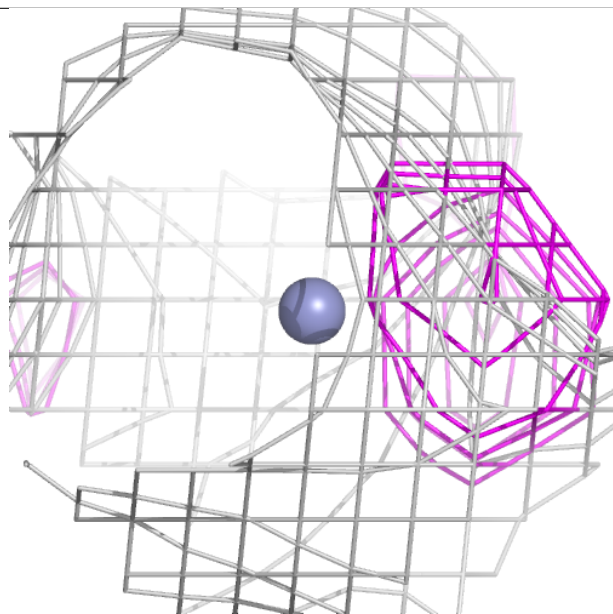
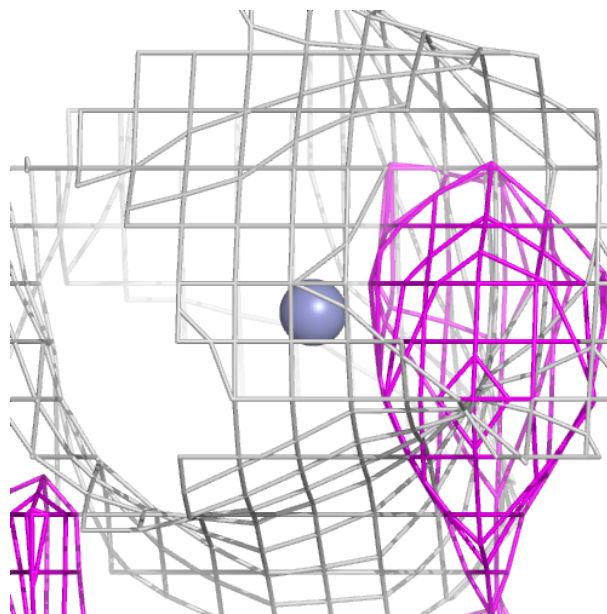
**Electron density around ZN D 401:**

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



**Electron density around ZN A 402:**

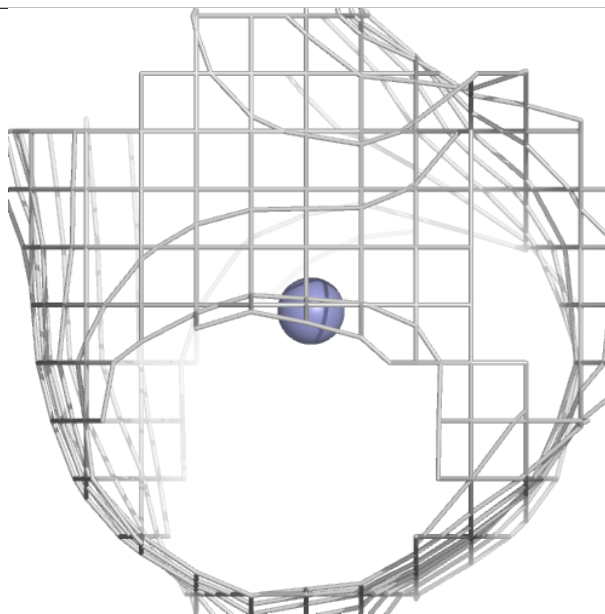
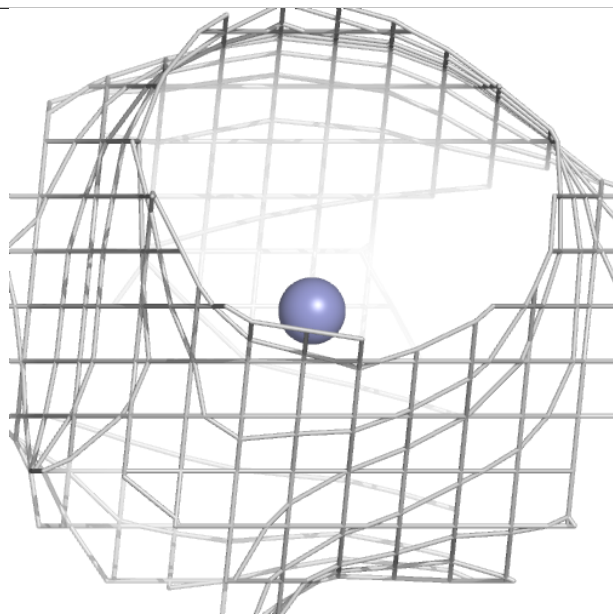
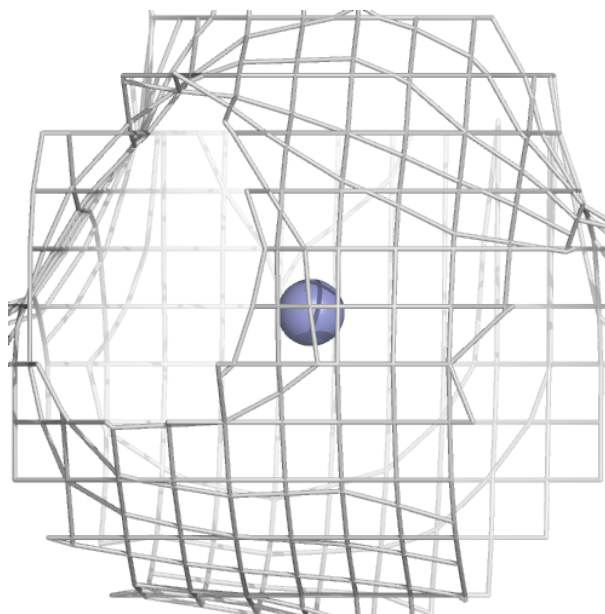
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





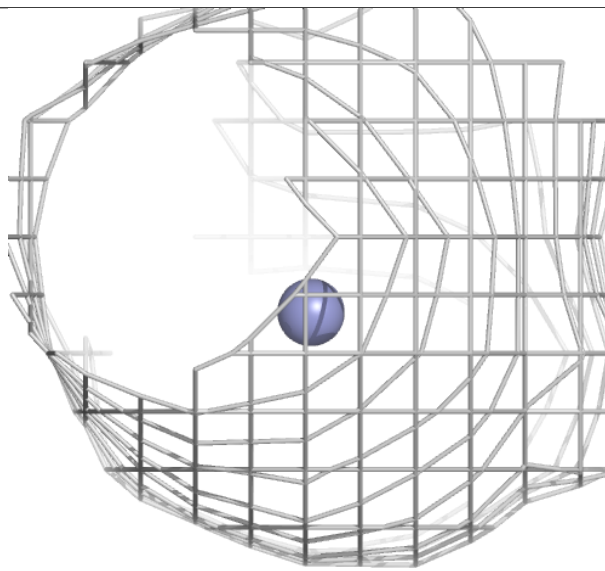
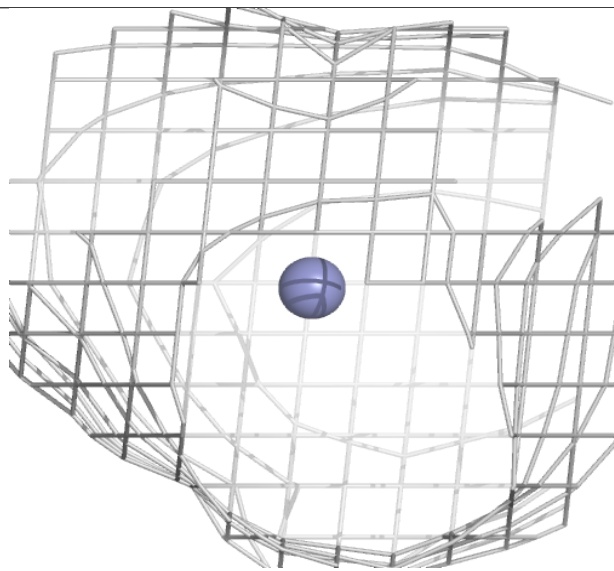
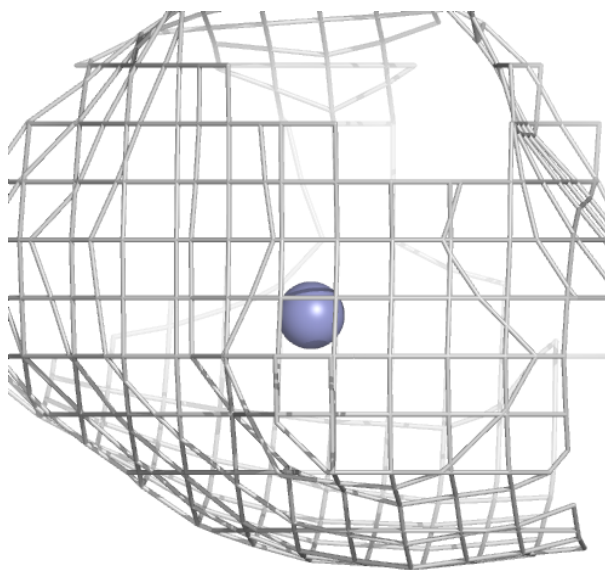
**Electron density around ZN B 401:**

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



**Electron density around ZN A 401:**

$2mF_o-DF_c$  (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.