



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2026 – 12:21 AM JST

PDB ID : 7W47 / pdb_00007w47
Title : Crystal structure of the gastric proton pump complexed with tegoprazan
Authors : Abe, K.; Tanaka, S.; Morita, M.; Yamagishi, T.
Deposited on : 2021-11-26
Resolution : 3.00 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 : 4-5-2 with Phenix2.0
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

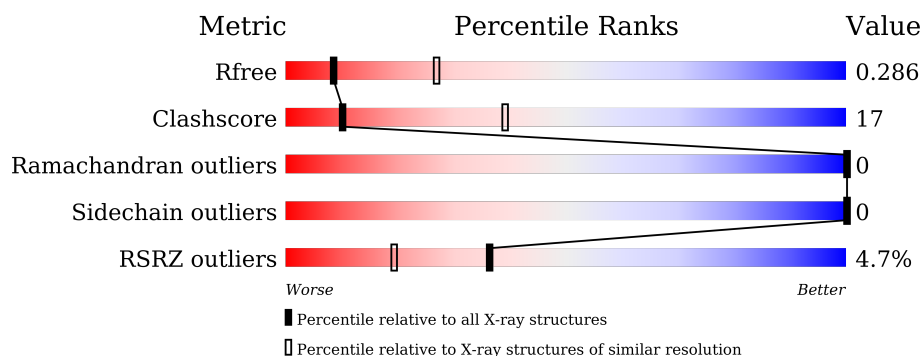
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	1034	
2	B	290	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9904 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 Potassium-transporting ATPase alpha chain 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	987	Total	C	N	O	S	0	0	0
			7659	4890	1292	1423	54			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	220	CYS	ARG	engineered mutation	UNP P19156
A	593	CYS	SER	engineered mutation	UNP P19156
A	1005	SER	GLY	engineered mutation	UNP P19156

- Molecule 2 is a protein called Potassium-transporting ATPase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	260	Total	C	N	O	S	0	0	0
			2084	1353	346	374	11			

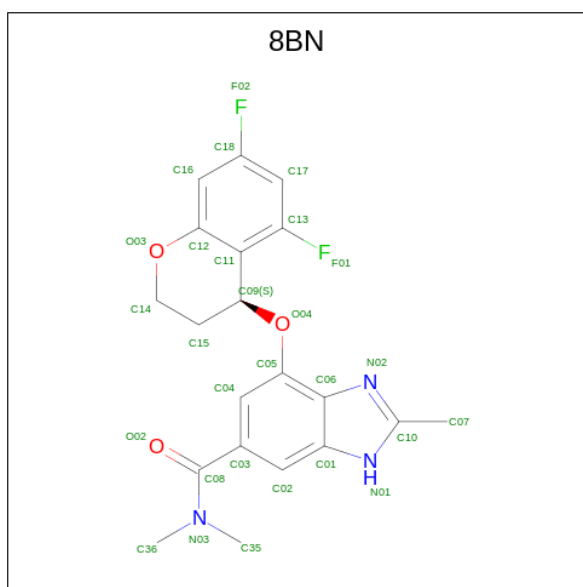
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is RUBIDIUM ION (CCD ID: RB) (formula: Rb) (labeled as "Ligand of Interest" by depositor).

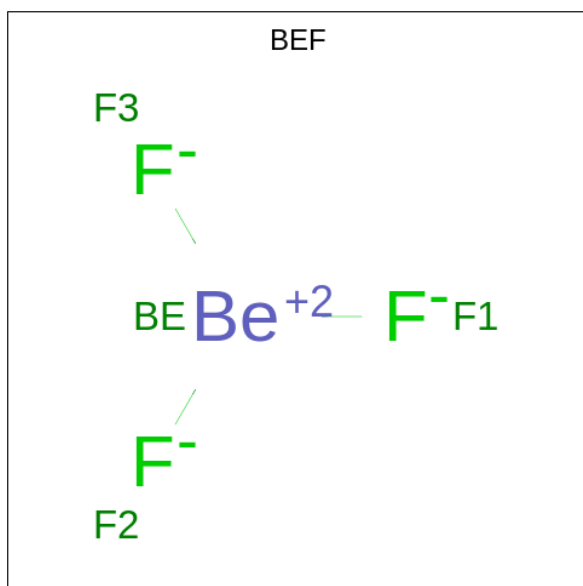
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Rb	0	0
			3	3		

- Molecule 5 is Tegoprazan (CCD ID: 8BN) (formula: C₂₀H₁₉F₂N₃O₃) (labeled as "Ligand of Interest" by depositor).



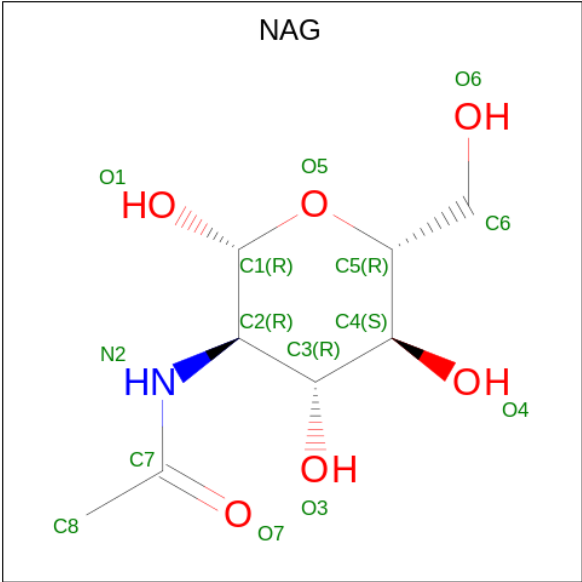
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	N	O	0
			28	20	2	3	3	

- Molecule 6 is BERYLLIUM TRIFLUORIDE ION (CCD ID: BEF) (formula: BeF₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Be	F	0	0
			4	1	3		

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		

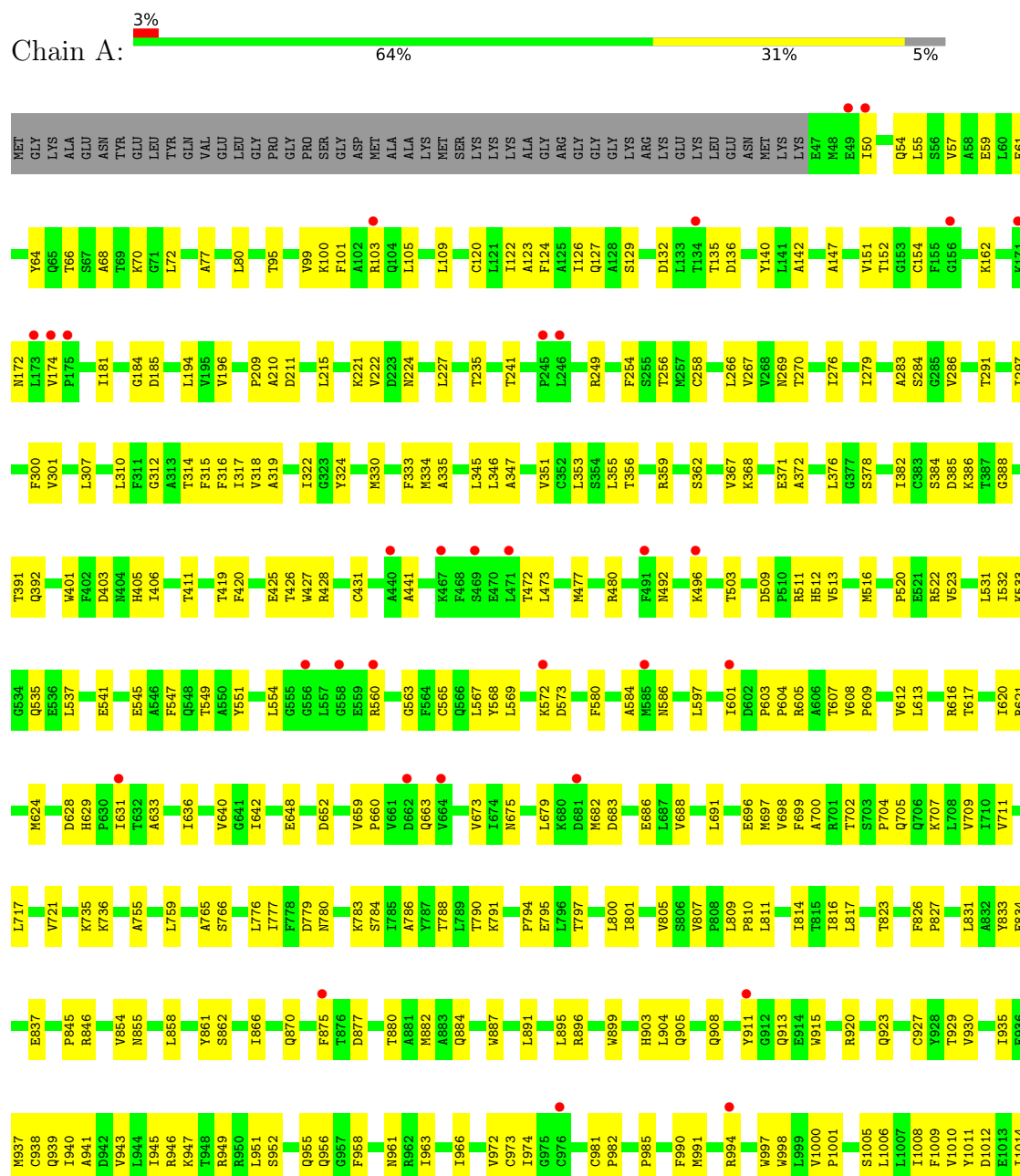
- Molecule 8 is water.

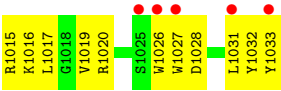
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	63	Total	O	0	0
			63	63		
8	B	20	Total	O	0	0
			20	20		

3 Residue-property plots

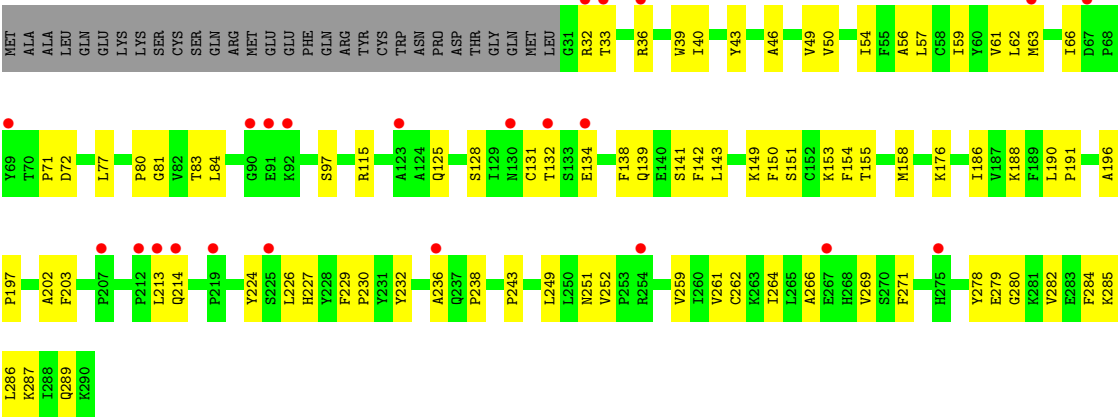
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

- Molecule 1: Potassium-transporting ATPase alpha chain 1





● Molecule 2: Potassium-transporting ATPase subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.99Å 104.99Å 367.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.26 – 3.00 48.26 – 3.00	Depositor EDS
% Data completeness (in resolution range)	84.8 (48.26-3.00) 84.7 (48.26-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.222 , 0.285 0.224 , 0.286	Depositor DCC
R_{free} test set	2028 reflections (2.31%)	wwPDB-VP
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 32.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.045 for -h,-k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	9904	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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: NAG, 8BN, BEF, MG, 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	0.48	1/7817 (0.0%)	0.74	3/10618 (0.0%)
2	B	0.43	0/2151	0.75	0/2926
All	All	0.47	1/9968 (0.0%)	0.74	3/13544 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	385	ASP	CG-OD2	9.26	1.43	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	384	SER	N-CA-C	5.60	117.53	108.41
1	A	973	CYS	CA-C-N	-5.53	110.79	120.29
1	A	973	CYS	C-N-CA	-5.53	110.79	120.29

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	7659	0	7684	267	0
2	B	2084	0	2025	82	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
4	A	3	0	0	0	0
5	A	28	0	0	1	0
6	A	4	0	0	0	0
7	B	42	0	39	1	0
8	A	63	0	0	9	0
8	B	20	0	0	5	0
All	All	9904	0	9748	337	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 17.

All (337) 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)
2:B:128:SER:HB2	2:B:153:LYS:HG3	1.44	0.99
2:B:139:GLN:HE22	2:B:150:PHE:HB2	1.27	0.96
1:A:441:ALA:HA	1:A:477:MET:HE1	1.50	0.93
1:A:882:MET:HE2	1:A:923:GLN:HE21	1.35	0.91
1:A:985:PRO:HG3	1:A:991:MET:HG2	1.56	0.86
1:A:929:THR:HG21	1:A:991:MET:H	1.43	0.82
2:B:214:GLN:HG2	2:B:251:ASN:HB2	1.61	0.80
2:B:141:SER:O	2:B:149:LYS:NZ	2.13	0.78
2:B:139:GLN:HB3	2:B:149:LYS:HE3	1.64	0.78
1:A:227:LEU:HD21	1:A:276:ILE:HD11	1.65	0.77
1:A:109:LEU:HB2	1:A:346:LEU:HD13	1.69	0.75
1:A:624:MET:HE1	1:A:633:ALA:HB1	1.69	0.75
1:A:945:ILE:HD11	1:A:1009:PHE:HA	1.69	0.74
2:B:128:SER:HB2	2:B:153:LYS:HZ2	1.53	0.74
2:B:139:GLN:NE2	2:B:150:PHE:HB2	2.02	0.74
1:A:222:VAL:HG23	1:A:258:CYS:HA	1.70	0.73
1:A:516:MET:HE3	1:A:523:VAL:HG23	1.69	0.73
1:A:1011:TYR:HH	2:B:43:TYR:HH	1.36	0.73
1:A:101:PHE:HE1	1:A:154:CYS:HB3	1.54	0.72
2:B:139:GLN:OE1	2:B:149:LYS:HD2	1.90	0.72
1:A:837:GLU:HB3	1:A:951:LEU:HD13	1.70	0.71
1:A:801:ILE:HG12	1:A:875:PHE:CZ	2.26	0.70
1:A:877:ASP:HB3	1:A:930:VAL:HG22	1.73	0.70
1:A:805:VAL:HG23	1:A:807:VAL:HG13	1.74	0.69
2:B:128:SER:HB2	2:B:153:LYS:CG	2.21	0.69
1:A:887:TRP:HH2	1:A:915:TRP:CD2	2.11	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:911:TYR:HE2	2:B:71:PRO:HA	1.58	0.68
1:A:699:PHE:HB3	1:A:702:THR:HG21	1.77	0.67
1:A:908:GLN:HA	1:A:913:GLN:O	1.93	0.67
1:A:545:GLU:O	1:A:549:THR:HG23	1.94	0.67
1:A:880:THR:HG22	1:A:997:TRP:HE1	1.58	0.67
1:A:391:THR:HA	1:A:604:PRO:HA	1.77	0.67
2:B:259:VAL:HG22	2:B:285:LYS:HG3	1.78	0.66
2:B:196:ALA:HA	2:B:226:LEU:HD21	1.77	0.66
1:A:882:MET:HE3	1:A:927:CYS:SG	2.36	0.65
1:A:801:ILE:HG12	1:A:875:PHE:HZ	1.61	0.65
1:A:431:CYS:SG	1:A:473:LEU:HD13	2.37	0.65
1:A:120:CYS:HB2	1:A:142:ALA:HB2	1.79	0.64
1:A:1016:LYS:HA	1:A:1019:VAL:HG12	1.78	0.64
1:A:605:ARG:HB2	1:A:608:VAL:HG23	1.80	0.64
1:A:1015:ARG:HH22	1:A:1032:TYR:HA	1.63	0.64
1:A:560:ARG:NH1	8:A:1203:HOH:O	2.30	0.63
1:A:1015:ARG:NH2	1:A:1032:TYR:HA	2.13	0.63
1:A:816:ILE:HG21	8:A:1228:HOH:O	1.99	0.63
2:B:46:ALA:HA	2:B:49:VAL:HG22	1.81	0.63
1:A:823:THR:HG22	1:A:974:ILE:HD11	1.81	0.63
2:B:32:ARG:HG2	2:B:40:ILE:HD11	1.81	0.62
1:A:783:LYS:HZ3	1:A:949:ARG:HG3	1.64	0.62
1:A:403:ASP:CG	1:A:426:THR:HG21	2.24	0.61
2:B:252:VAL:HG21	2:B:286:LEU:HD23	1.82	0.61
1:A:834:GLU:OE2	1:A:947:LYS:HE3	2.01	0.61
1:A:172:ASN:HD21	1:A:174:VAL:HG12	1.65	0.61
1:A:795:GLU:HB2	8:A:1228:HOH:O	2.01	0.61
1:A:356:THR:OG1	1:A:777:ILE:HG21	2.01	0.61
2:B:289:GLN:HB2	7:B:301:NAG:H82	1.83	0.61
1:A:532:ILE:HG22	1:A:533:LYS:HG2	1.82	0.60
1:A:998:TRP:O	1:A:1001:PRO:HD2	2.01	0.60
2:B:128:SER:CB	2:B:153:LYS:HG3	2.23	0.60
1:A:624:MET:CE	1:A:698:VAL:HG13	2.31	0.60
1:A:403:ASP:OD1	1:A:426:THR:HG21	2.02	0.60
1:A:994:ARG:HD2	8:B:418:HOH:O	2.01	0.60
1:A:794:PRO:HB2	1:A:935:ILE:HD11	1.84	0.59
1:A:1019:VAL:HG23	1:A:1028:ASP:OD1	2.03	0.59
1:A:378:SER:OG	1:A:846:ARG:NH1	2.35	0.59
1:A:927:CYS:HA	1:A:930:VAL:HG12	1.84	0.59
2:B:226:LEU:HD11	8:B:416:HOH:O	2.02	0.58
2:B:271:PHE:HA	2:B:280:GLY:HA3	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:TRP:HB3	1:A:597:LEU:HB2	1.85	0.58
1:A:963:ILE:HA	1:A:966:ILE:HG22	1.85	0.58
1:A:227:LEU:HD11	1:A:276:ILE:HD13	1.85	0.58
1:A:291:THR:HG21	1:A:371:GLU:HB3	1.85	0.58
1:A:345:LEU:HD13	1:A:788:THR:HG21	1.85	0.58
1:A:427:TRP:HZ2	1:A:472:THR:HG21	1.69	0.58
1:A:613:LEU:O	1:A:617:THR:HG23	2.03	0.58
1:A:707:LYS:O	1:A:711:VAL:HG23	2.03	0.58
2:B:128:SER:HB3	2:B:151:SER:O	2.03	0.58
2:B:197:PRO:HG2	2:B:224:TYR:HB2	1.86	0.58
1:A:315:PHE:CE2	1:A:800:LEU:HB2	2.39	0.58
1:A:425:GLU:HG2	1:A:428:ARG:HH21	1.67	0.58
1:A:861:TYR:CE2	1:A:866:ILE:HD11	2.39	0.58
2:B:81:GLY:HA2	2:B:279:GLU:HA	1.85	0.57
1:A:833:TYR:HB2	1:A:961:ASN:HD21	1.69	0.57
1:A:127:GLN:HB3	1:A:132:ASP:HB3	1.86	0.57
1:A:210:ALA:O	1:A:254:PHE:HA	2.04	0.57
1:A:899:TRP:O	1:A:920:ARG:NH1	2.37	0.57
2:B:264:ILE:HG22	2:B:269:VAL:HG21	1.86	0.57
1:A:705:GLN:O	1:A:709:VAL:HG23	2.05	0.56
1:A:887:TRP:CH2	1:A:915:TRP:CE2	2.94	0.56
2:B:132:THR:O	2:B:134:GLU:HG3	2.05	0.56
1:A:1012:ASP:OD1	1:A:1015:ARG:NE	2.36	0.56
1:A:388:GLY:O	1:A:605:ARG:NH1	2.39	0.56
1:A:958:PHE:HZ	1:A:1009:PHE:CZ	2.24	0.56
1:A:624:MET:HE2	1:A:698:VAL:HG13	1.88	0.55
1:A:347:ALA:O	1:A:351:VAL:HG22	2.06	0.55
1:A:57:VAL:O	1:A:61:GLU:HG2	2.06	0.55
1:A:427:TRP:CZ2	1:A:472:THR:HG21	2.41	0.55
1:A:858:LEU:HB2	1:A:1033:TYR:HB2	1.89	0.55
1:A:1017:LEU:HD12	8:A:1223:HOH:O	2.07	0.55
1:A:68:ALA:HB1	1:A:215:LEU:HD13	1.89	0.54
1:A:985:PRO:CG	1:A:991:MET:HG2	2.31	0.54
1:A:181:ILE:HA	1:A:185:ASP:O	2.07	0.54
2:B:46:ALA:O	2:B:50:VAL:HG23	2.08	0.54
2:B:128:SER:H	2:B:153:LYS:HZ3	1.56	0.54
1:A:1026:TRP:HA	2:B:36:ARG:HH22	1.73	0.54
1:A:120:CYS:CB	1:A:142:ALA:HB2	2.38	0.53
1:A:382:ILE:HG13	1:A:620:ILE:HG21	1.90	0.53
1:A:814:ILE:HG22	8:A:1227:HOH:O	2.07	0.53
1:A:1017:LEU:HG	1:A:1020:ARG:HH21	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:MET:HE2	1:A:565:CYS:SG	2.48	0.53
1:A:319:ALA:O	1:A:324:TYR:HB2	2.08	0.53
1:A:905:GLN:HB2	2:B:83:THR:HG22	1.91	0.53
1:A:496:LYS:HB3	1:A:522:ARG:NE	2.24	0.53
1:A:759:LEU:HD22	1:A:766:SER:HB2	1.89	0.53
2:B:188:LYS:HA	2:B:230:PRO:HB2	1.91	0.53
1:A:794:PRO:CG	1:A:870:GLN:HB2	2.38	0.53
1:A:941:ALA:CB	1:A:1008:ILE:HD11	2.38	0.53
1:A:503:THR:HG22	1:A:512:HIS:CD2	2.44	0.53
1:A:866:ILE:HD12	1:A:866:ILE:H	1.74	0.53
1:A:866:ILE:HD12	1:A:866:ILE:N	2.24	0.53
1:A:95:THR:OG1	1:A:100:LYS:HD2	2.09	0.52
1:A:673:VAL:HG13	1:A:698:VAL:HG12	1.90	0.52
1:A:952:SER:HB3	1:A:955:GLN:HG3	1.92	0.52
2:B:50:VAL:O	2:B:54:ILE:HG13	2.08	0.52
1:A:866:ILE:O	1:A:870:GLN:HG3	2.10	0.52
1:A:101:PHE:CE1	1:A:154:CYS:HB3	2.42	0.52
1:A:659:VAL:HG22	1:A:660:PRO:HD2	1.91	0.52
1:A:531:LEU:O	1:A:532:ILE:HD13	2.10	0.52
1:A:783:LYS:NZ	1:A:949:ARG:HG3	2.24	0.52
1:A:915:TRP:CE2	2:B:77:LEU:HD21	2.45	0.52
2:B:36:ARG:O	2:B:40:ILE:HG13	2.10	0.52
1:A:963:ILE:HD12	1:A:963:ILE:H	1.75	0.52
2:B:80:PRO:HB3	2:B:186:ILE:HD11	1.91	0.52
1:A:511:ARG:HB2	1:A:569:LEU:O	2.10	0.51
1:A:862:SER:HB3	1:A:946:ARG:HH22	1.76	0.51
1:A:911:TYR:CE2	2:B:71:PRO:HA	2.42	0.51
1:A:791:LYS:HD2	1:A:939:GLN:NE2	2.26	0.51
1:A:563:GLY:HA2	1:A:597:LEU:HD23	1.93	0.51
1:A:941:ALA:HB1	1:A:1008:ILE:HD11	1.91	0.51
2:B:33:THR:OG1	2:B:36:ARG:HB2	2.09	0.51
1:A:318:VAL:O	1:A:322:ILE:HG12	2.11	0.51
1:A:882:MET:CE	1:A:923:GLN:HE21	2.17	0.51
1:A:711:VAL:HG13	1:A:721:VAL:HG11	1.93	0.51
2:B:32:ARG:HA	2:B:32:ARG:NE	2.26	0.51
1:A:122:ILE:HG13	1:A:334:MET:HE1	1.92	0.51
1:A:368:LYS:HD2	1:A:755:ALA:O	2.11	0.51
1:A:816:ILE:HG13	1:A:817:LEU:N	2.26	0.51
2:B:203:PHE:CZ	2:B:213:LEU:HB3	2.46	0.51
1:A:794:PRO:HB2	1:A:935:ILE:CD1	2.40	0.50
1:A:884:GLN:HG3	2:B:72:ASP:HB2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:929:THR:CG2	1:A:990:PHE:HA	2.41	0.50
1:A:127:GLN:HE21	1:A:135:THR:HG22	1.76	0.50
1:A:406:ILE:HD13	1:A:554:LEU:HD21	1.93	0.50
1:A:784:SER:O	1:A:788:THR:HG22	2.12	0.50
1:A:324:TYR:OH	1:A:896:ARG:NH2	2.44	0.50
1:A:831:LEU:HD11	1:A:943:VAL:HG23	1.93	0.50
2:B:202:ALA:HA	2:B:213:LEU:HD11	1.94	0.50
1:A:50:ILE:HD12	1:A:50:ILE:H	1.77	0.49
1:A:783:LYS:NZ	1:A:837:GLU:OE1	2.33	0.49
1:A:392:GLN:HG3	1:A:603:PRO:O	2.13	0.49
1:A:826:PHE:HB3	1:A:827:PRO:HD3	1.95	0.49
1:A:64:TYR:CE1	1:A:196:VAL:HG12	2.47	0.49
1:A:1015:ARG:HG3	1:A:1031:LEU:HD23	1.93	0.49
1:A:70:LYS:HE3	1:A:184:GLY:HA3	1.95	0.49
1:A:241:THR:HB	1:A:249:ARG:H	1.77	0.49
1:A:648:GLU:HB2	1:A:652:ASP:HB2	1.94	0.49
1:A:605:ARG:HB2	1:A:608:VAL:CG2	2.43	0.49
1:A:386:LYS:NZ	1:A:633:ALA:HA	2.28	0.49
1:A:316:PHE:HB2	1:A:333:PHE:HB2	1.95	0.49
1:A:1005:SER:O	1:A:1008:ILE:HG13	2.13	0.49
1:A:310:LEU:C	1:A:310:LEU:HD23	2.38	0.48
1:A:172:ASN:ND2	1:A:174:VAL:HG12	2.28	0.48
1:A:580:PHE:HA	1:A:586:ASN:HD21	1.78	0.48
1:A:887:TRP:HH2	1:A:915:TRP:CE2	2.31	0.48
2:B:66:ILE:HD12	2:B:66:ILE:O	2.14	0.48
1:A:861:TYR:CZ	1:A:866:ILE:HD11	2.48	0.48
1:A:122:ILE:HG13	1:A:334:MET:CE	2.44	0.48
1:A:124:PHE:CD2	1:A:135:THR:HB	2.49	0.48
1:A:547:PHE:CD1	1:A:547:PHE:C	2.92	0.48
1:A:688:VAL:HG12	1:A:717:LEU:HD11	1.94	0.48
1:A:880:THR:HG22	1:A:997:TRP:NE1	2.28	0.48
1:A:659:VAL:HG22	1:A:663:GLN:HG3	1.95	0.48
1:A:679:LEU:HD12	1:A:682:MET:HE2	1.94	0.48
1:A:683:ASP:OD2	1:A:686:GLU:HG3	2.13	0.48
2:B:149:LYS:HA	2:B:149:LYS:HD3	1.52	0.48
1:A:862:SER:HB3	1:A:946:ARG:NH2	2.30	0.47
1:A:310:LEU:O	1:A:314:THR:HG23	2.15	0.47
1:A:797:THR:HA	1:A:800:LEU:HD11	1.95	0.47
1:A:72:LEU:HD21	1:A:80:LEU:HD12	1.96	0.47
2:B:57:LEU:O	2:B:61:VAL:HG23	2.15	0.47
2:B:202:ALA:HA	2:B:213:LEU:CD1	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:GLU:HB2	1:A:66:THR:O	2.14	0.47
1:A:683:ASP:OD2	1:A:686:GLU:N	2.38	0.47
1:A:279:ILE:HD13	1:A:704:PRO:HG2	1.97	0.47
1:A:297:ILE:HA	1:A:300:PHE:CE2	2.50	0.47
2:B:261:VAL:HA	2:B:282:VAL:O	2.15	0.47
1:A:221:LYS:CD	1:A:235:THR:HG22	2.45	0.47
2:B:59:ILE:HA	2:B:62:LEU:HB3	1.97	0.47
1:A:55:LEU:HD22	1:A:59:GLU:HB3	1.97	0.47
1:A:72:LEU:HD23	1:A:77:ALA:HA	1.95	0.47
1:A:1015:ARG:HA	1:A:1027:TRP:HZ2	1.79	0.47
2:B:36:ARG:O	2:B:39:TRP:N	2.48	0.47
2:B:176:LYS:HA	2:B:249:LEU:O	2.15	0.47
1:A:227:LEU:HD21	1:A:276:ILE:CD1	2.40	0.46
1:A:367:VAL:CG2	1:A:372:ALA:HB3	2.46	0.46
1:A:284:SER:HA	1:A:735:LYS:HE3	1.96	0.46
1:A:1026:TRP:HA	2:B:36:ARG:NH2	2.29	0.46
2:B:97:SER:HA	2:B:287:LYS:O	2.16	0.46
1:A:503:THR:HG22	1:A:512:HIS:NE2	2.31	0.46
1:A:162:LYS:HD2	1:A:162:LYS:HA	1.64	0.46
2:B:142:PHE:HA	2:B:149:LYS:HE2	1.98	0.46
1:A:609:PRO:O	1:A:613:LEU:HG	2.15	0.46
2:B:285:LYS:N	2:B:285:LYS:HD2	2.31	0.46
1:A:101:PHE:CE2	1:A:105:LEU:HD11	2.50	0.46
1:A:126:ILE:O	1:A:129:SER:HB3	2.16	0.46
1:A:887:TRP:CH2	1:A:915:TRP:CD2	2.99	0.46
2:B:125:GLN:OE1	2:B:154:PHE:N	2.47	0.46
1:A:330:MET:HB3	1:A:330:MET:HE2	1.48	0.45
1:A:779:ASP:O	1:A:783:LYS:HG3	2.17	0.45
1:A:795:GLU:HA	1:A:811:LEU:HD23	1.98	0.45
1:A:887:TRP:CD1	1:A:891:LEU:HD23	2.51	0.45
1:A:903:HIS:HE1	8:B:419:HOH:O	1.99	0.45
1:A:359:ARG:O	1:A:362:SER:OG	2.33	0.45
2:B:131:CYS:HB3	2:B:227:HIS:CE1	2.51	0.45
1:A:572:LYS:HD3	1:A:572:LYS:C	2.42	0.45
2:B:32:ARG:HE	2:B:33:THR:H	1.63	0.45
2:B:128:SER:N	2:B:153:LYS:HZ3	2.15	0.45
1:A:477:MET:HE2	1:A:477:MET:HB3	1.86	0.45
1:A:777:ILE:HA	1:A:780:ASN:HB2	1.98	0.45
2:B:142:PHE:HD2	2:B:149:LYS:HG2	1.81	0.45
2:B:262:CYS:HB2	2:B:282:VAL:CG2	2.47	0.45
1:A:1017:LEU:HG	1:A:1020:ARG:NH2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:LEU:CD1	1:A:209:PRO:HB2	2.46	0.44
1:A:572:LYS:HD3	1:A:573:ASP:N	2.32	0.44
1:A:604:PRO:HG3	1:A:636:ILE:HD12	1.99	0.44
1:A:607:THR:HG22	8:A:1220:HOH:O	2.18	0.44
1:A:951:LEU:HD22	1:A:956:GLN:NE2	2.32	0.44
1:A:317:ILE:HG13	1:A:318:VAL:N	2.31	0.44
1:A:631:ILE:HD12	1:A:631:ILE:H	1.83	0.44
1:A:855:ASN:OD1	1:A:855:ASN:N	2.50	0.44
1:A:136:ASP:HB3	1:A:140:TYR:CE2	2.52	0.44
1:A:147:ALA:O	1:A:151:VAL:HG23	2.17	0.44
1:A:675:ASN:HA	1:A:700:ALA:O	2.17	0.44
1:A:981:CYS:SG	1:A:982:PRO:HD2	2.58	0.44
1:A:1010:VAL:O	1:A:1014:ILE:HG12	2.16	0.44
1:A:441:ALA:HA	1:A:477:MET:CE	2.33	0.44
1:A:776:LEU:HD23	1:A:776:LEU:HA	1.79	0.44
2:B:138:PHE:HZ	2:B:190:LEU:HD21	1.82	0.44
2:B:224:TYR:HD1	2:B:243:PRO:HG2	1.83	0.44
1:A:109:LEU:HD23	1:A:301:VAL:CG1	2.47	0.44
1:A:152:THR:HG21	1:A:346:LEU:HD23	2.00	0.44
1:A:211:ASP:HB2	1:A:269:ASN:O	2.17	0.44
1:A:411:THR:HB	1:A:603:PRO:CB	2.48	0.44
1:A:790:THR:HG22	1:A:870:GLN:OE1	2.16	0.44
1:A:958:PHE:HZ	1:A:1009:PHE:HZ	1.66	0.44
2:B:63:MET:HA	2:B:66:ILE:HD11	1.98	0.44
1:A:535:GLN:HB2	1:A:537:LEU:CD1	2.47	0.43
2:B:236:ALA:O	2:B:238:PRO:HD3	2.18	0.43
1:A:99:VAL:O	1:A:103:ARG:HG2	2.19	0.43
1:A:858:LEU:O	1:A:862:SER:OG	2.28	0.43
1:A:427:TRP:CH2	1:A:431:CYS:HB2	2.53	0.43
1:A:628:ASP:OD1	1:A:629:HIS:N	2.50	0.43
1:A:513:VAL:HB	1:A:568:TYR:CE2	2.54	0.43
1:A:640:VAL:HG23	1:A:642:ILE:HG12	2.00	0.43
1:A:845:PRO:HA	8:A:1211:HOH:O	2.17	0.43
1:A:736:LYS:HB2	1:A:736:LYS:HE2	1.80	0.43
1:A:891:LEU:HG	1:A:895:LEU:HD13	2.01	0.43
1:A:794:PRO:HG3	1:A:870:GLN:HB2	2.01	0.43
1:A:642:ILE:O	1:A:696:GLU:HB3	2.18	0.43
2:B:155:THR:H	2:B:158:MET:HE3	1.84	0.43
1:A:520:PRO:HG2	1:A:551:TYR:CE1	2.54	0.42
1:A:904:LEU:C	1:A:905:GLN:HG2	2.44	0.42
1:A:283:ALA:O	1:A:286:VAL:HG22	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:PHE:HB3	1:A:854:VAL:HG22	2.01	0.42
1:A:312:GLY:C	1:A:333:PHE:CD2	2.98	0.42
1:A:297:ILE:HG21	1:A:297:ILE:HD13	1.70	0.42
1:A:477:MET:HA	1:A:480:ARG:HG2	2.00	0.42
1:A:854:VAL:HA	1:A:858:LEU:HD23	2.01	0.42
1:A:401:TRP:HA	1:A:405:HIS:O	2.19	0.42
1:A:509:ASP:OD1	1:A:511:ARG:HG2	2.19	0.42
2:B:224:TYR:HE1	2:B:243:PRO:O	2.03	0.42
2:B:266:ALA:HB3	2:B:269:VAL:HG22	2.01	0.42
1:A:267:VAL:HG21	1:A:270:THR:HG22	2.01	0.42
1:A:441:ALA:CA	1:A:477:MET:HE1	2.35	0.42
1:A:621:ARG:NH2	1:A:691:LEU:O	2.41	0.42
1:A:817:LEU:HD23	1:A:817:LEU:HA	1.80	0.42
2:B:197:PRO:HB3	2:B:266:ALA:HB2	2.01	0.42
1:A:224:ASN:OD1	1:A:256:THR:HG21	2.20	0.42
1:A:937:MET:HE2	1:A:937:MET:HA	2.02	0.42
2:B:191:PRO:HG3	2:B:229:PHE:CE2	2.53	0.42
2:B:66:ILE:HG22	2:B:72:ASP:CG	2.44	0.42
1:A:376:LEU:HA	1:A:376:LEU:HD23	1.65	0.42
1:A:419:THR:OG1	1:A:420:PHE:N	2.51	0.42
1:A:607:THR:OG1	1:A:765:ALA:HB2	2.20	0.42
1:A:215:LEU:HD11	1:A:266:LEU:HG	2.02	0.41
1:A:1006:LEU:O	1:A:1010:VAL:HG23	2.20	0.41
1:A:307:LEU:HD23	1:A:307:LEU:HA	1.68	0.41
1:A:567:LEU:HD22	1:A:569:LEU:HD23	2.02	0.41
1:A:697:MET:HE2	1:A:697:MET:HB3	1.74	0.41
1:A:797:THR:HA	1:A:800:LEU:CD1	2.50	0.41
2:B:32:ARG:HG3	2:B:36:ARG:HB3	2.01	0.41
2:B:66:ILE:HG22	2:B:72:ASP:OD2	2.20	0.41
2:B:142:PHE:CD2	2:B:149:LYS:HG2	2.56	0.41
1:A:353:LEU:HD23	1:A:777:ILE:HD12	2.02	0.41
1:A:427:TRP:HZ2	1:A:472:THR:CG2	2.31	0.41
2:B:143:LEU:O	2:B:150:PHE:CZ	2.73	0.41
1:A:516:MET:CE	1:A:523:VAL:HG23	2.45	0.41
1:A:997:TRP:HA	1:A:1000:VAL:HG23	2.02	0.41
1:A:492:ASN:OD1	1:A:492:ASN:N	2.52	0.41
1:A:809:LEU:HD12	1:A:810:PRO:HD2	2.03	0.41
1:A:335:ALA:HB2	5:A:1105:8BN:C13	2.51	0.41
1:A:791:LYS:C	1:A:794:PRO:HD2	2.45	0.41
2:B:56:ALA:O	2:B:59:ILE:HG13	2.20	0.41
1:A:123:ALA:HB2	1:A:334:MET:HE2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:541:GLU:HG3	8:A:1258:HOH:O	2.21	0.41
1:A:584:ALA:O	1:A:586:ASN:N	2.54	0.41
1:A:905:GLN:HG3	2:B:278:TYR:CD2	2.56	0.41
1:A:940:ILE:HD11	1:A:972:VAL:CG1	2.51	0.41
1:A:601:ILE:HD12	1:A:601:ILE:HA	1.87	0.41
1:A:904:LEU:O	1:A:920:ARG:NH2	2.54	0.41
1:A:1016:LYS:HA	1:A:1019:VAL:CG1	2.49	0.41
2:B:203:PHE:HD2	8:B:408:HOH:O	2.04	0.41
2:B:230:PRO:HG2	2:B:232:TYR:HE1	1.85	0.41
1:A:425:GLU:HG2	1:A:428:ARG:NH2	2.35	0.41
1:A:612:VAL:O	1:A:616:ARG:HG3	2.21	0.41
1:A:816:ILE:HD11	8:A:1263:HOH:O	2.21	0.41
1:A:905:GLN:HG3	2:B:278:TYR:CE2	2.56	0.41
1:A:809:LEU:HD12	1:A:810:PRO:CD	2.51	0.40
1:A:884:GLN:HG3	2:B:72:ASP:CB	2.51	0.40
1:A:786:ALA:HB2	1:A:858:LEU:HD11	2.03	0.40
1:A:880:THR:HA	2:B:62:LEU:HD11	2.02	0.40
2:B:84:LEU:HD11	2:B:284:PHE:CE1	2.57	0.40
2:B:115:ARG:HD3	2:B:115:ARG:HA	1.92	0.40
1:A:870:GLN:HE21	1:A:938:CYS:HB2	1.85	0.40
2:B:115:ARG:NH2	8:B:401:HOH:O	2.13	0.40
1:A:50:ILE:O	1:A:54:GLN:NE2	2.47	0.40
1:A:351:VAL:O	1:A:355:LEU:HG	2.20	0.40
1:A:346:LEU:HD12	1:A:346:LEU:HA	1.91	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	985/1034 (95%)	935 (95%)	50 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	258/290 (89%)	239 (93%)	19 (7%)	0	100	100
All	All	1243/1324 (94%)	1174 (94%)	69 (6%)	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	834/870 (96%)	834 (100%)	0	100	100
2	B	227/254 (89%)	227 (100%)	0	100	100
All	All	1061/1124 (94%)	1061 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	GLN
1	A	422	GLN
1	A	542	GLN
1	A	629	HIS
1	A	773	GLN
1	A	870	GLN
1	A	898	GLN
1	A	923	GLN
1	A	986	ASN
2	B	76	GLN
2	B	139	GLN
2	B	273	ASN
2	B	289	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 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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$
7	NAG	B	303	2	14,14,15	0.55	0	17,19,21	0.90	0
7	NAG	B	302	2	14,14,15	0.28	0	17,19,21	0.52	0
5	8BN	A	1105	-	31,31,31	2.87	7 (22%)	36,46,46	2.49	14 (38%)
7	NAG	B	301	2	14,14,15	0.57	1 (7%)	17,19,21	1.16	2 (11%)
6	BEF	A	1106	1	0,3,3	-	-	-	-	-

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
7	NAG	B	303	2	-	1/6/23/26	0/1/1/1
7	NAG	B	302	2	-	2/6/23/26	0/1/1/1
5	8BN	A	1105	-	-	1/12/22/22	0/4/4/4
7	NAG	B	301	2	-	2/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1105	8BN	C11-C13	11.01	1.50	1.38
5	A	1105	8BN	C01-C06	7.37	1.47	1.40
5	A	1105	8BN	C12-C11	5.04	1.47	1.39
5	A	1105	8BN	C05-C06	4.28	1.47	1.40
5	A	1105	8BN	C06-N02	-2.90	1.34	1.39
5	A	1105	8BN	C08-N03	-2.44	1.31	1.34
5	A	1105	8BN	C01-N01	-2.24	1.34	1.38
7	B	301	NAG	C1-C2	2.04	1.55	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1105	8BN	C14-O03-C12	6.39	124.68	113.65
5	A	1105	8BN	F01-C13-C11	6.06	124.06	118.13
5	A	1105	8BN	C17-C13-C11	-4.73	118.69	124.63
5	A	1105	8BN	C13-C17-C18	4.31	121.14	116.62
5	A	1105	8BN	C05-C06-N02	4.03	132.72	126.55
5	A	1105	8BN	O03-C12-C16	3.28	122.10	117.05
7	B	301	NAG	C1-O5-C5	3.27	116.62	112.19
5	A	1105	8BN	C16-C12-C11	-3.27	116.96	121.85
5	A	1105	8BN	C12-C16-C18	2.92	122.39	116.45
5	A	1105	8BN	C07-C10-N01	2.67	126.93	122.78
5	A	1105	8BN	O04-C09-C15	2.38	112.85	108.49
5	A	1105	8BN	C03-C04-C05	2.32	122.70	119.61
5	A	1105	8BN	C02-C01-N01	2.22	135.26	130.65
7	B	301	NAG	C1-C2-N2	2.08	114.05	110.49
5	A	1105	8BN	O04-C09-C11	-2.07	102.54	106.76
5	A	1105	8BN	C04-C05-C06	-2.03	116.78	120.45

There are no chirality outliers.

All (6) torsion outliers are listed below:

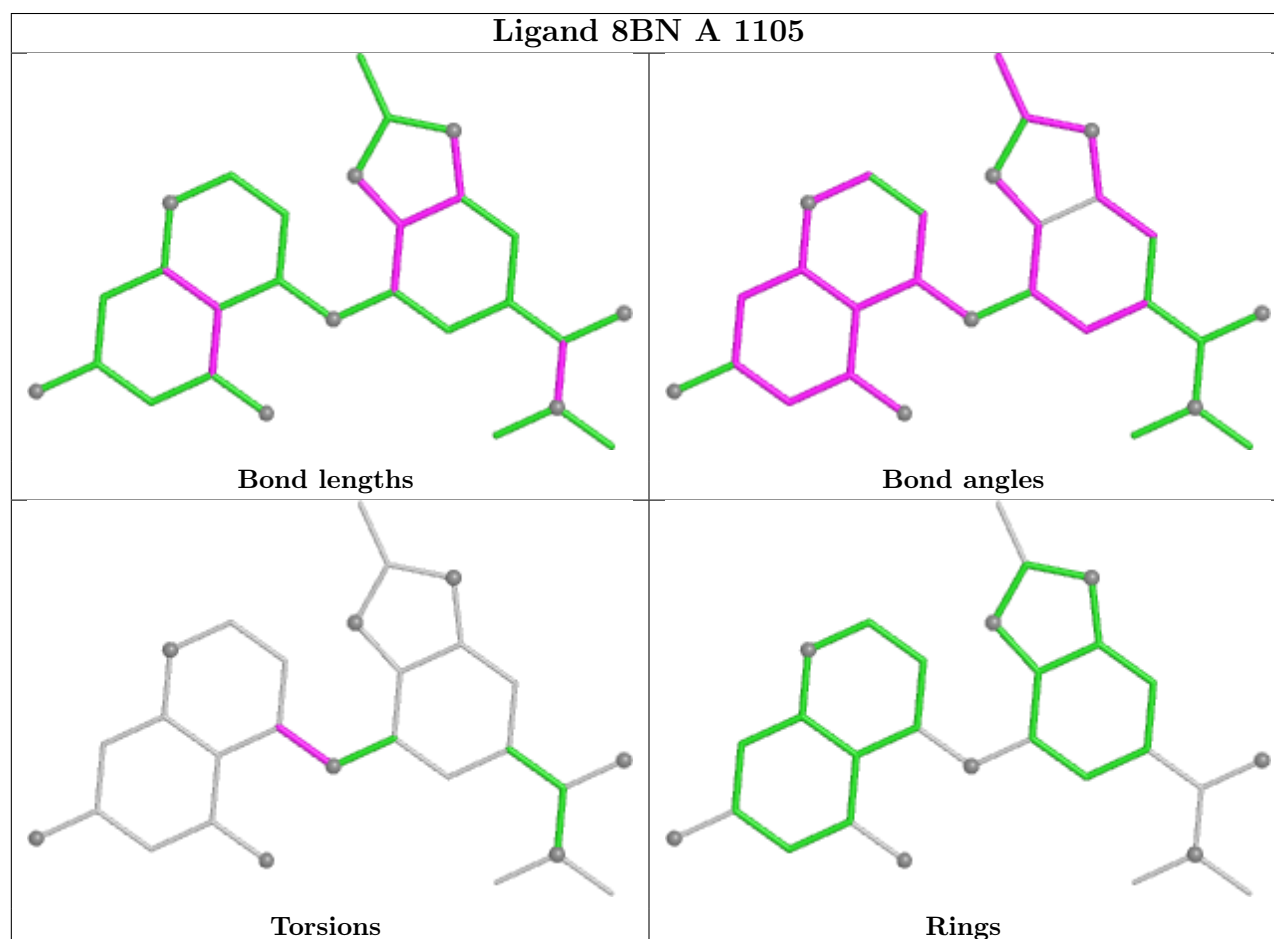
Mol	Chain	Res	Type	Atoms
7	B	302	NAG	C4-C5-C6-O6
7	B	302	NAG	O5-C5-C6-O6
5	A	1105	8BN	C11-C09-O04-C05
7	B	301	NAG	O5-C5-C6-O6
7	B	301	NAG	C4-C5-C6-O6
7	B	303	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1105	8BN	1	0
7	B	301	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	987/1034 (95%)	0.09	36 (3%)	46 26	22, 52, 99, 166	0
2	B	260/290 (89%)	0.55	23 (8%)	15 8	36, 73, 124, 191	0
All	All	1247/1324 (94%)	0.19	59 (4%)	36 19	22, 57, 108, 191	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	33	THR	6.4
2	B	130	ASN	5.6
1	A	662	ASP	5.4
2	B	132	THR	5.2
1	A	50	ILE	4.7
1	A	49	GLU	4.3
2	B	92	LYS	4.2
2	B	91	GLU	4.0
2	B	90	GLY	4.0
2	B	236	ALA	4.0
1	A	558	GLY	3.9
1	A	246	LEU	3.9
2	B	212	PRO	3.9
2	B	267	GLU	3.8
2	B	69	TYR	3.5
1	A	1033	TYR	3.3
1	A	469	SER	3.2
2	B	32	ARG	3.2
1	A	171	LYS	3.2
2	B	67	ASP	3.1
1	A	601	ILE	3.1
1	A	585	MET	3.0
1	A	1027	TRP	2.9
1	A	1025	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	976	CYS	2.8
2	B	63	MET	2.8
1	A	994	ARG	2.7
2	B	134	GLU	2.7
1	A	491	PHE	2.6
1	A	496	LYS	2.6
1	A	572	LYS	2.6
2	B	213	LEU	2.5
1	A	467	LYS	2.4
1	A	174	VAL	2.4
1	A	631	ILE	2.4
1	A	245	PRO	2.4
1	A	1026	TRP	2.4
1	A	1031	LEU	2.3
2	B	225	SER	2.3
1	A	556	GLY	2.2
2	B	275	HIS	2.2
2	B	207	PRO	2.2
1	A	173	LEU	2.2
1	A	134	THR	2.2
1	A	471	LEU	2.2
1	A	103	ARG	2.2
2	B	36	ARG	2.2
2	B	214	GLN	2.2
2	B	219	PRO	2.2
1	A	681	ASP	2.2
1	A	664	VAL	2.1
2	B	254	ARG	2.1
1	A	156	GLY	2.1
1	A	911	TYR	2.1
1	A	560	ARG	2.1
1	A	440	ALA	2.1
1	A	875	PHE	2.1
2	B	123	ALA	2.1
1	A	175	PRO	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

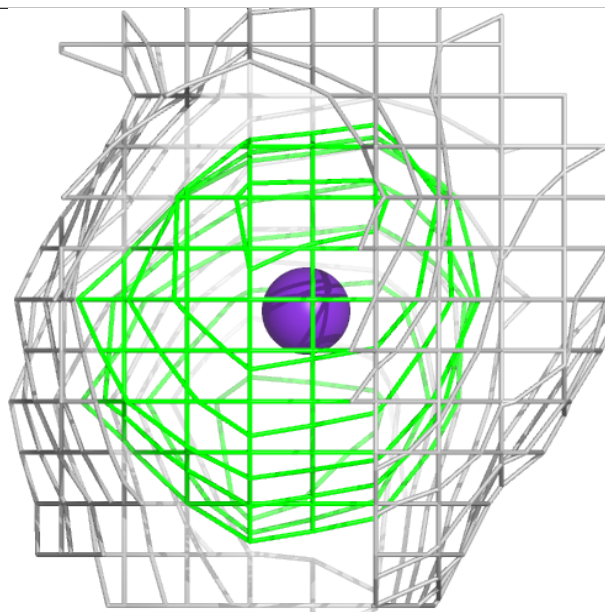
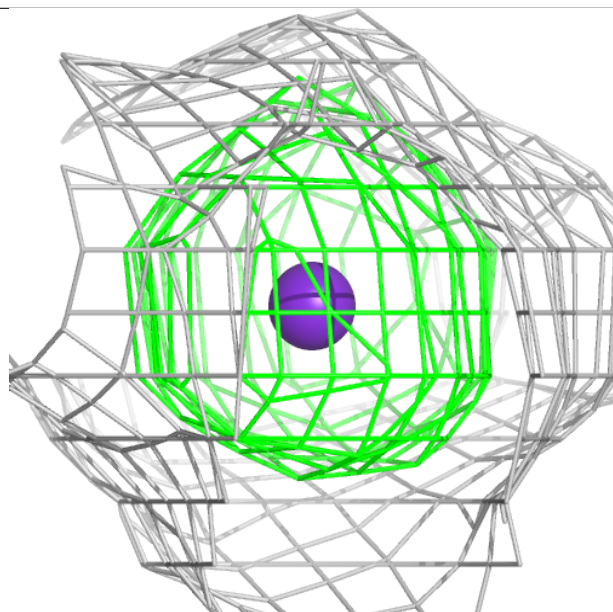
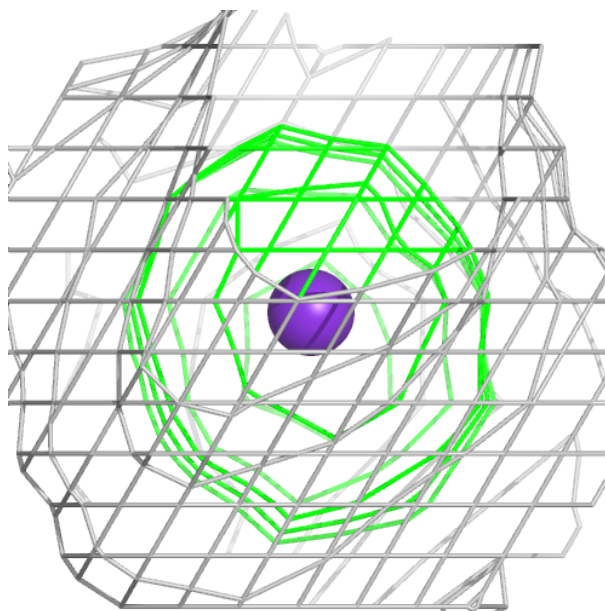
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
7	NAG	B	302	14/15	0.73	0.22	110,122,131,132	0
7	NAG	B	303	14/15	0.88	0.15	86,112,121,121	0
7	NAG	B	301	14/15	0.90	0.10	34,48,53,64	0
6	BEF	A	1106	4/4	0.94	0.14	38,40,42,43	0
4	RB	A	1104	1/1	0.95	0.20	130,130,130,130	0
5	8BN	A	1105	28/28	0.95	0.10	25,43,68,78	0
3	MG	A	1101	1/1	0.98	0.13	21,21,21,21	0
4	RB	A	1102	1/1	0.99	0.11	95,95,95,95	0
4	RB	A	1103	1/1	0.99	0.08	73,73,73,73	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.

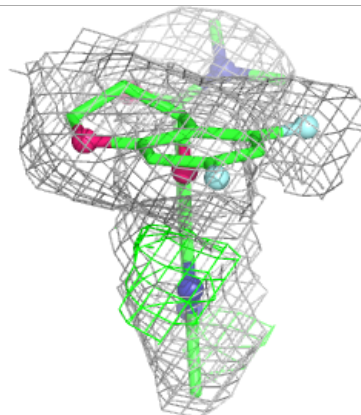
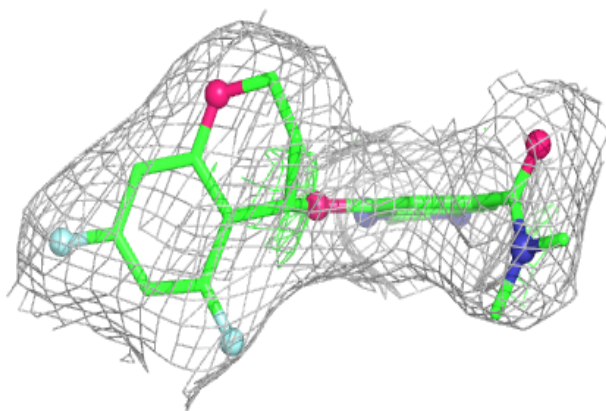
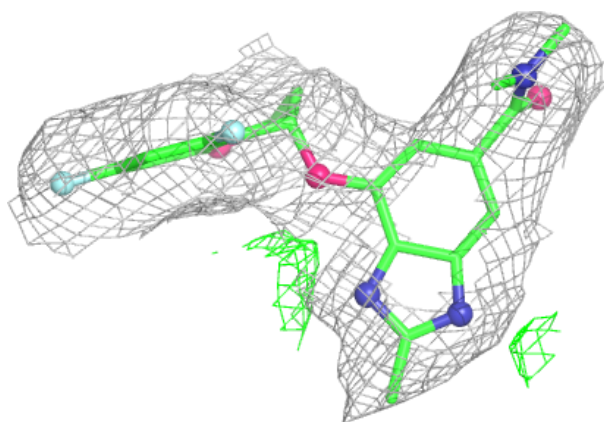
Electron density around RB A 1104:

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



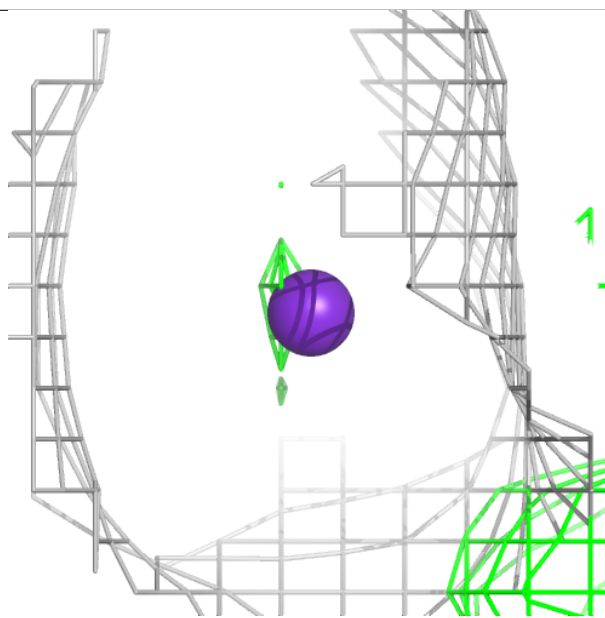
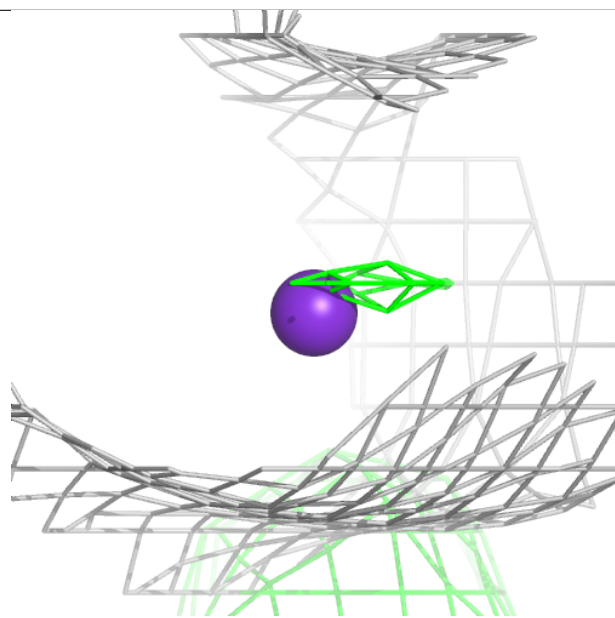
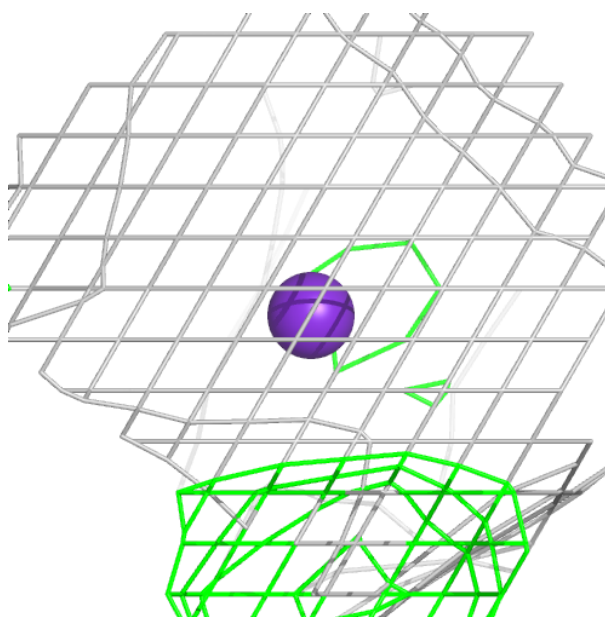
Electron density around 8BN A 1105:

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



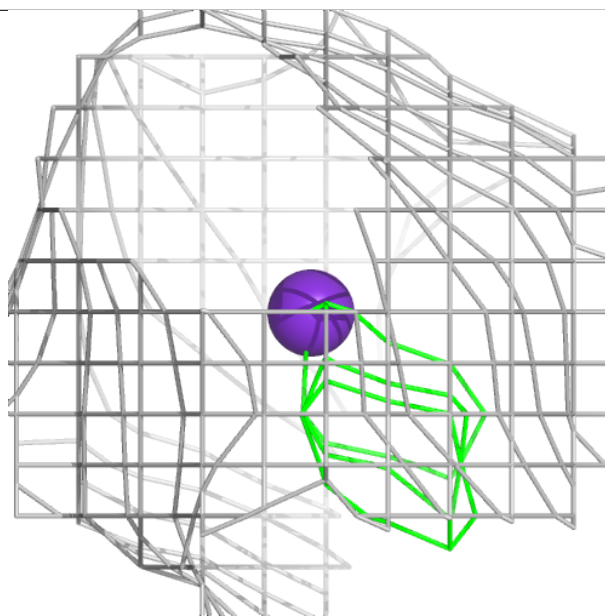
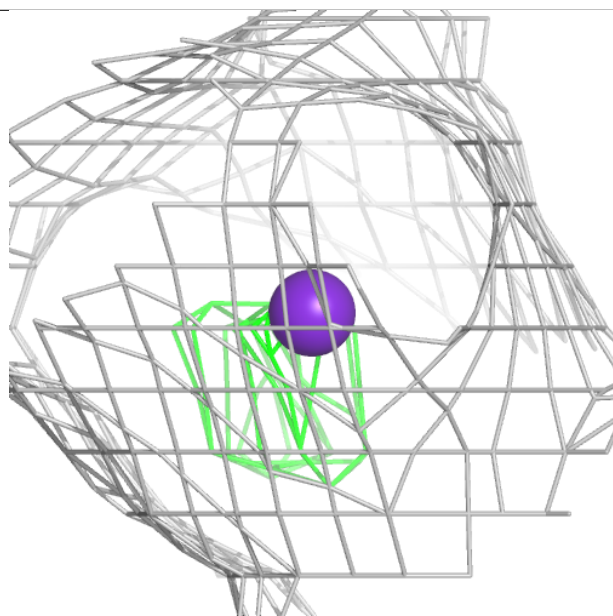
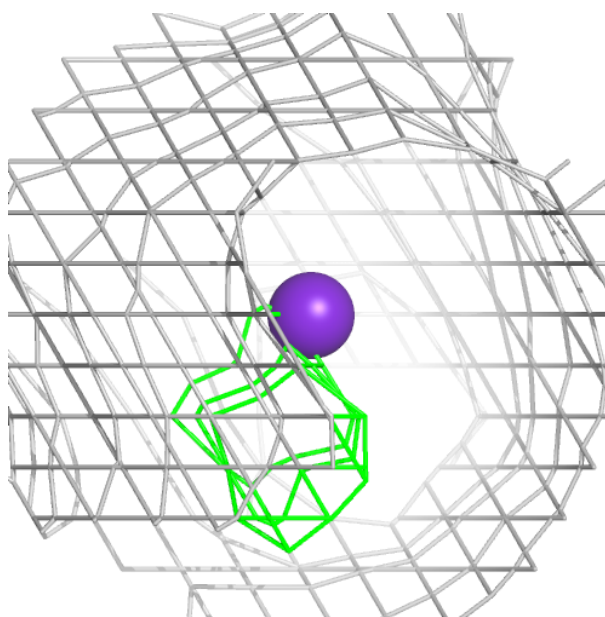
Electron density around RB A 1102:

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



Electron density around RB A 1103:

$2mF_o-DF_c$ (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.