



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:29 AM UTC

PDB ID : 7FKT / pdb_00007fkt
Title : PanDDA analysis group deposition – Aar2/RNaseH in complex with fragment P04F02 from the F2X-Universal Library
Authors : Barthel, T.; Wollenhaupt, J.; Lima, G.M.A.; Wahl, M.C.; Weiss, M.S.
Deposited on : 2022-08-26
Resolution : 1.44 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

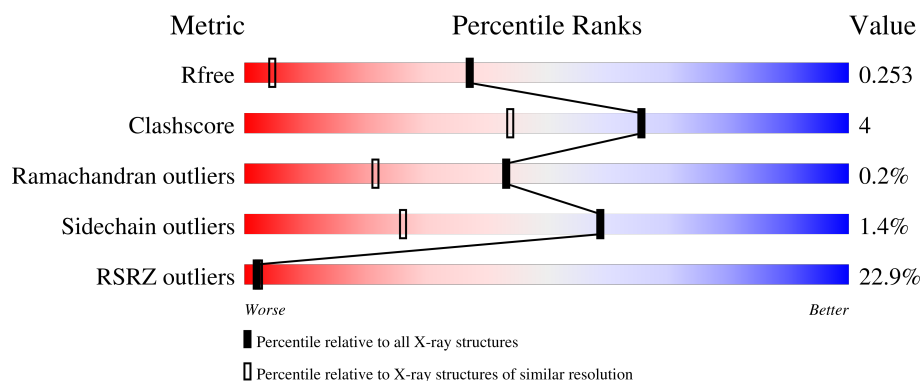
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.44 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	3234 (1.46-1.42)
Clashscore	190562	3289 (1.46-1.42)
Ramachandran outliers	187476	3248 (1.46-1.42)
Sidechain outliers	187428	3248 (1.46-1.42)
RSRZ outliers	180081	3234 (1.46-1.42)

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	258	18% 85% 7% 8%
2	B	308	25% 86% 10% ••

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
3	VBI	B	401	-	X	-	-
3	VBI	B	402	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9308 atoms, of which 4515 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 Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	237	4068	1287	2060	336	373	12	0	21	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	GLY	-	expression tag	UNP P33334
A	1834	ALA	-	expression tag	UNP P33334
A	1835	MET	-	expression tag	UNP P33334

- Molecule 2 is a protein called A1 cistron-splicing factor AAR2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	300	5027	1649	2455	420	484	19	11	16	0

There are 20 discrepancies between the modelled and reference sequences:

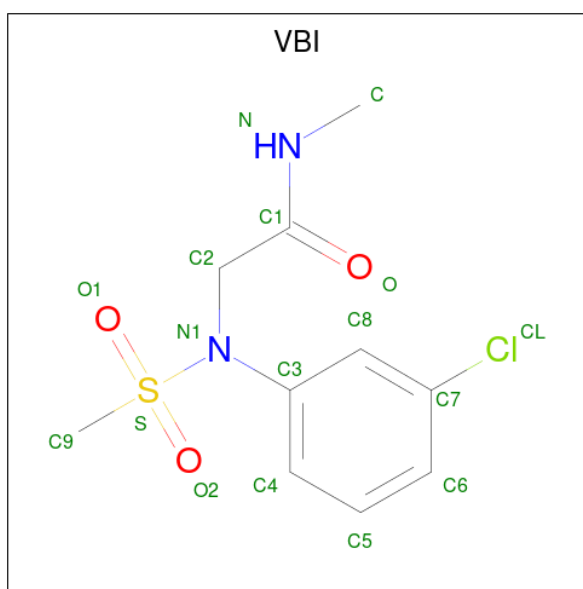
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	LEU	conflict	UNP P32357
B	167	SER	LYS	conflict	UNP P32357
B	?	-	LEU	deletion	UNP P32357
B	?	-	GLN	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	GLY	deletion	UNP P32357
B	?	-	SER	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	MET	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ASN	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	170	SER	ASP	conflict	UNP P32357

- Molecule 3 is N 2 -(3-chlorophenyl)-N 2 -(methanesulfonyl)-N-methylglycinamide (CCD ID: VBI) (formula: C₁₀H₁₃ClN₂O₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			17	10	1	2	3	1		
3	B	1	Total	C	Cl	N	O	S	0	0
			17	10	1	2	3	1		
3	B	1	Total	C	Cl	N	O	S	0	0
			17	10	1	2	3	1		
3	B	1	Total	C	Cl	N	O	S	0	0
			17	10	1	2	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	77	Total	O	0	0
			77	77		

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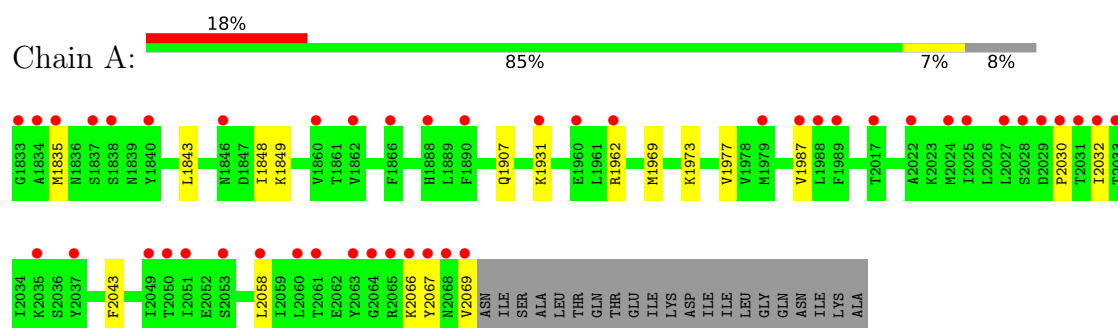
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	68	Total	O	0	0
			68	68		

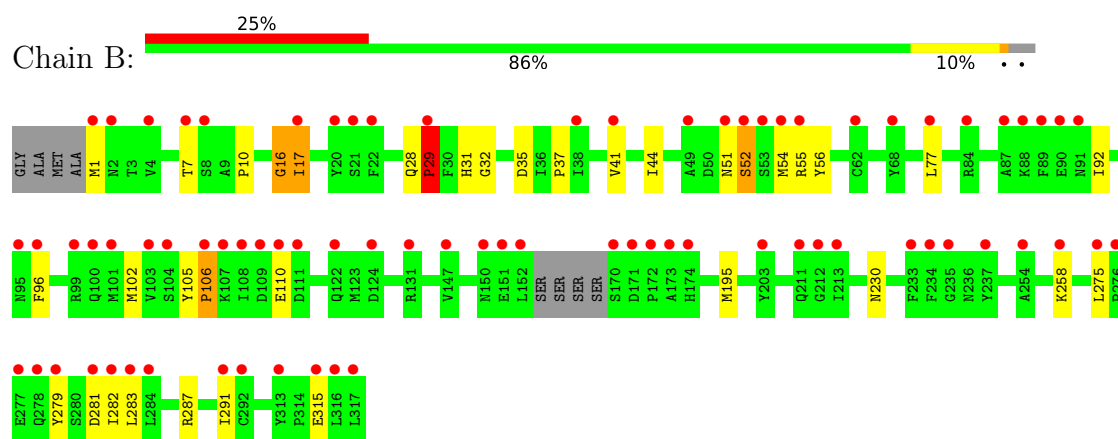
3 Residue-property plots [i](#)

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: Pre-mRNA-splicing factor 8



- Molecule 2: A1 cistron-splicing factor AAR2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	89.27Å 81.78Å 93.78Å 90.00° 108.57° 90.00°	Depositor
Resolution (Å)	44.45 – 1.44 44.45 – 1.44	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.45-1.44) 99.9 (44.45-1.44)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 1.44Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.210 , 0.247 0.231 , 0.253	Depositor DCC
R_{free} test set	2099 reflections (1.82%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9308	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% 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: VBI

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	0.53	0/2149	0.72	3/2911 (0.1%)
2	B	0.74	4/2731 (0.1%)	1.01	16/3689 (0.4%)
All	All	0.66	4/4880 (0.1%)	0.90	19/6600 (0.3%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	29	PRO	N-CA	18.08	1.70	1.47
2	B	230[A]	ASN	C-O	7.51	1.32	1.24
2	B	230[B]	ASN	C-O	7.51	1.32	1.24
2	B	28	GLN	C-N	5.08	1.45	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	28	GLN	CA-C-N	13.16	136.29	119.84
2	B	28	GLN	C-N-CA	13.16	136.29	119.84
2	B	230[A]	ASN	CA-C-O	10.11	131.27	120.55
2	B	230[B]	ASN	CA-C-O	10.11	131.27	120.55
2	B	29	PRO	CA-N-CD	-8.91	99.53	112.00
2	B	106	PRO	N-CA-CB	-6.33	98.57	102.81
2	B	230[A]	ASN	N-CA-C	6.31	118.15	111.28
2	B	230[B]	ASN	N-CA-C	6.31	118.15	111.28
2	B	32	GLY	N-CA-C	6.19	119.32	110.46
1	A	1973	LYS	CA-C-O	-6.14	114.04	120.55
2	B	28	GLN	O-C-N	-5.76	114.97	121.43
2	B	230[A]	ASN	O-C-N	-5.55	116.24	122.12
2	B	230[B]	ASN	O-C-N	-5.55	116.24	122.12
2	B	10	PRO	CB-CA-C	-5.40	103.80	111.68
2	B	17	ILE	N-CA-C	-5.25	98.09	107.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	16	GLY	O-C-N	-5.17	117.01	123.21
2	B	31	HIS	CA-CB-CG	-5.06	108.74	113.80
1	A	2030	PRO	CA-C-N	5.05	127.56	120.28
1	A	2030	PRO	C-N-CA	5.05	127.56	120.28

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	2008	2060	1974	11	2
2	B	2572	2455	2389	25	2
3	A	17	0	0	0	0
3	B	51	0	0	0	0
4	A	77	0	0	0	0
4	B	68	0	0	3	0
All	All	4793	4515	4363	35	2

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 4.

All (35) 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:29:PRO:N	2:B:29:PRO:CA	1.70	1.41
2:B:275:LEU:CD2	2:B:283:LEU:HD13	2.09	0.83
2:B:275:LEU:HD22	2:B:283:LEU:CD1	2.12	0.79
2:B:275:LEU:HD22	2:B:283:LEU:HD12	1.69	0.74
2:B:275:LEU:HD21	2:B:283:LEU:HD13	1.70	0.72
2:B:275:LEU:CD2	2:B:283:LEU:CD1	2.72	0.67
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.43	0.65
2:B:1:MET:HB3	2:B:35:ASP:HA	1.86	0.57
2:B:37:PRO:HD3	2:B:105:TYR:CD1	2.41	0.56
2:B:51:ASN:ND2	4:B:502:HOH:O	2.38	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:258:LYS:HD2	2:B:258:LYS:H	1.71	0.55
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.33	0.54
2:B:17:ILE:O	2:B:17:ILE:HG23	2.11	0.51
2:B:1:MET:N	4:B:503:HOH:O	2.44	0.49
2:B:29:PRO:N	2:B:29:PRO:C	2.61	0.49
2:B:17:ILE:HD13	2:B:44[B]:ILE:HG12	1.95	0.47
2:B:56:TYR:CZ	2:B:77:LEU:HB2	2.49	0.47
2:B:96:PHE:CB	2:B:102:MET:HE3	2.44	0.47
2:B:17:ILE:HD13	2:B:44[B]:ILE:CG1	2.45	0.47
1:A:1969:MET:HA	1:A:1969:MET:HE2	1.97	0.47
1:A:2032:ILE:HG21	1:A:2043:PHE:CD1	2.51	0.46
1:A:1977:VAL:HG21	1:A:1987:VAL:HG21	1.97	0.45
2:B:287:ARG:O	2:B:291:ILE:HD13	2.17	0.45
2:B:17:ILE:CD1	2:B:44[B]:ILE:CG1	2.95	0.44
2:B:315:GLU:H	2:B:315:GLU:CD	2.25	0.43
1:A:1843:LEU:HA	1:A:1849:LYS:HD2	2.00	0.43
1:A:2066:LYS:HD2	1:A:2067:TYR:CE1	2.53	0.43
2:B:41[A]:VAL:HG12	4:B:546:HOH:O	2.18	0.43
1:A:1962:ARG:HD3	1:A:1962:ARG:H	1.84	0.42
2:B:110:GLU:HG2	2:B:110:GLU:O	2.19	0.42
1:A:1907:GLN:HG2	2:B:195:MET:HE3	2.01	0.41
1:A:1835:MET:HE3	1:A:1835:MET:HB3	1.96	0.41
2:B:281:ASP:OD1	2:B:281:ASP:N	2.54	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2066:LYS:HG3	2:B:279:TYR:HE2[2_655]	1.34	0.26
1:A:2069:VAL:O	2:B:52:SER:OG[2_655]	2.15	0.05

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	258/258 (100%)	254 (98%)	4 (2%)	0	100	100
2	B	314/308 (102%)	306 (98%)	7 (2%)	1 (0%)	36	17
All	All	572/566 (101%)	560 (98%)	11 (2%)	1 (0%)	43	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	29	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	237/233 (102%)	237 (100%)	0	100	100
2	B	293/284 (103%)	285 (97%)	8 (3%)	39	8
All	All	530/517 (102%)	522 (98%)	8 (2%)	59	24

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

Mol	Chain	Res	Type
2	B	7	THR
2	B	52	SER
2	B	54	MET
2	B	55[A]	ARG
2	B	55[B]	ARG
2	B	92	ILE
2	B	106	PRO
2	B	282	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	51	ASN
2	B	70	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	VBI	B	401	-	17,17,17	4.57	11 (64%)	23,24,24	2.33	8 (34%)
3	VBI	B	403	-	17,17,17	3.03	4 (23%)	23,24,24	1.68	6 (26%)
3	VBI	B	402	-	17,17,17	2.43	5 (29%)	23,24,24	2.41	8 (34%)
3	VBI	A	2101	-	17,17,17	2.46	5 (29%)	23,24,24	1.51	5 (21%)

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
3	VBI	B	401	-	-	5/16/16/16	0/1/1/1
3	VBI	B	403	-	-	1/16/16/16	0/1/1/1
3	VBI	B	402	-	-	11/16/16/16	0/1/1/1
3	VBI	A	2101	-	-	8/16/16/16	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	VBI	S-N1	-9.55	1.54	1.64
3	B	401	VBI	C9-S	-8.81	1.57	1.75
3	B	403	VBI	O2-S	-7.76	1.32	1.43
3	B	401	VBI	O2-S	-6.98	1.33	1.43
3	B	402	VBI	O2-S	-6.42	1.34	1.43
3	B	403	VBI	C9-S	-6.37	1.62	1.75
3	B	401	VBI	O1-S	-5.46	1.35	1.43
3	B	402	VBI	C9-S	-5.25	1.64	1.75
3	A	2101	VBI	O2-S	-5.24	1.35	1.43
3	A	2101	VBI	C9-S	-5.11	1.65	1.75
3	B	401	VBI	C4-C3	5.09	1.48	1.39
3	B	401	VBI	C3-N1	-4.74	1.38	1.44
3	B	401	VBI	C2-N1	-4.65	1.40	1.48
3	A	2101	VBI	O1-S	-4.25	1.37	1.43
3	B	403	VBI	O1-S	-4.13	1.37	1.43
3	B	403	VBI	S-N1	-3.95	1.60	1.64
3	B	402	VBI	O1-S	-3.19	1.38	1.43
3	B	401	VBI	C8-C7	2.93	1.43	1.38
3	A	2101	VBI	C3-N1	2.90	1.48	1.44
3	B	401	VBI	C7-CL	-2.89	1.67	1.74
3	B	401	VBI	C5-C6	-2.84	1.34	1.38
3	A	2101	VBI	O-C1	-2.75	1.17	1.23
3	B	402	VBI	O-C1	-2.59	1.18	1.23
3	B	401	VBI	C6-C7	-2.40	1.33	1.38
3	B	402	VBI	C2-N1	-2.39	1.44	1.48

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	VBI	C6-C7-CL	-6.40	109.91	119.36
3	B	402	VBI	O1-S-C9	6.30	116.94	108.30
3	B	402	VBI	O2-S-N1	4.52	112.79	107.39
3	B	401	VBI	C8-C7-CL	4.22	124.46	119.17
3	B	403	VBI	C4-C3-N1	4.04	126.03	120.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	VBI	C9-S-N1	-3.86	100.31	106.41
3	B	403	VBI	C8-C7-CL	-3.78	114.43	119.17
3	B	402	VBI	O2-S-C9	-3.73	103.19	108.30
3	A	2101	VBI	C6-C7-CL	-3.48	114.23	119.36
3	B	401	VBI	C3-C8-C7	-3.39	114.36	118.92
3	B	402	VBI	C2-N1-S	3.33	123.53	118.09
3	B	401	VBI	C6-C7-C8	3.02	125.48	121.53
3	B	401	VBI	C4-C3-N1	3.00	124.53	120.15
3	B	402	VBI	C2-N1-C3	-2.97	112.68	116.67
3	B	401	VBI	C1-C2-N1	-2.96	106.97	115.43
3	B	402	VBI	O1-S-N1	-2.66	104.21	107.39
3	A	2101	VBI	C8-C3-N1	2.62	124.38	119.35
3	B	403	VBI	C8-C3-N1	-2.52	114.50	119.35
3	A	2101	VBI	C8-C7-CL	2.46	122.25	119.17
3	B	403	VBI	C2-N1-C3	2.36	119.83	116.67
3	B	401	VBI	O2-S-N1	-2.32	104.62	107.39
3	A	2101	VBI	C4-C3-N1	-2.28	116.82	120.15
3	B	401	VBI	O2-S-O1	2.04	121.50	118.46
3	B	403	VBI	O1-S-C9	2.02	111.08	108.30
3	B	403	VBI	O2-S-C9	-2.02	105.53	108.30
3	A	2101	VBI	C9-S-N1	-2.01	103.23	106.41
3	B	402	VBI	C6-C7-CL	-2.01	116.40	119.36

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2101	VBI	C3-N1-S-C9
3	A	2101	VBI	C3-N1-S-O1
3	A	2101	VBI	C2-N1-S-C9
3	A	2101	VBI	C2-N1-S-O1
3	A	2101	VBI	C2-C1-N-C
3	A	2101	VBI	O-C1-N-C
3	B	401	VBI	C3-N1-S-C9
3	B	401	VBI	C3-N1-S-O1
3	B	401	VBI	C2-N1-S-C9
3	B	401	VBI	C2-N1-S-O1
3	B	402	VBI	C8-C3-N1-C2
3	B	402	VBI	C4-C3-N1-C2
3	B	402	VBI	C8-C3-N1-S
3	B	402	VBI	C4-C3-N1-S
3	B	402	VBI	C3-N1-S-O1

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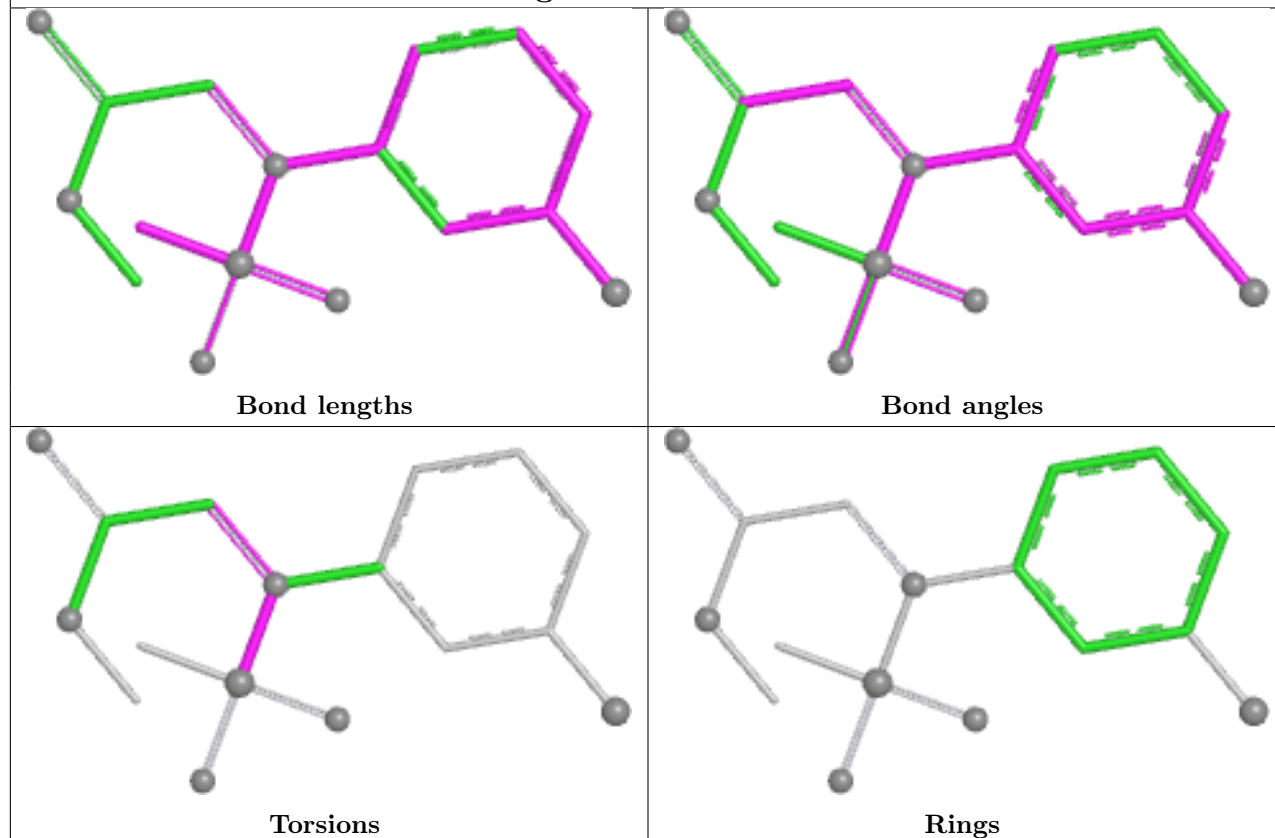
Mol	Chain	Res	Type	Atoms
3	B	402	VBI	C2-N1-S-O1
3	B	402	VBI	C2-N1-S-O2
3	B	402	VBI	C2-C1-N-C
3	B	402	VBI	O-C1-N-C
3	B	401	VBI	C1-C2-N1-S
3	B	402	VBI	C2-N1-S-C9
3	A	2101	VBI	O-C1-C2-N1
3	A	2101	VBI	N-C1-C2-N1
3	B	403	VBI	C1-C2-N1-C3
3	B	402	VBI	C3-N1-S-C9

There are no ring outliers.

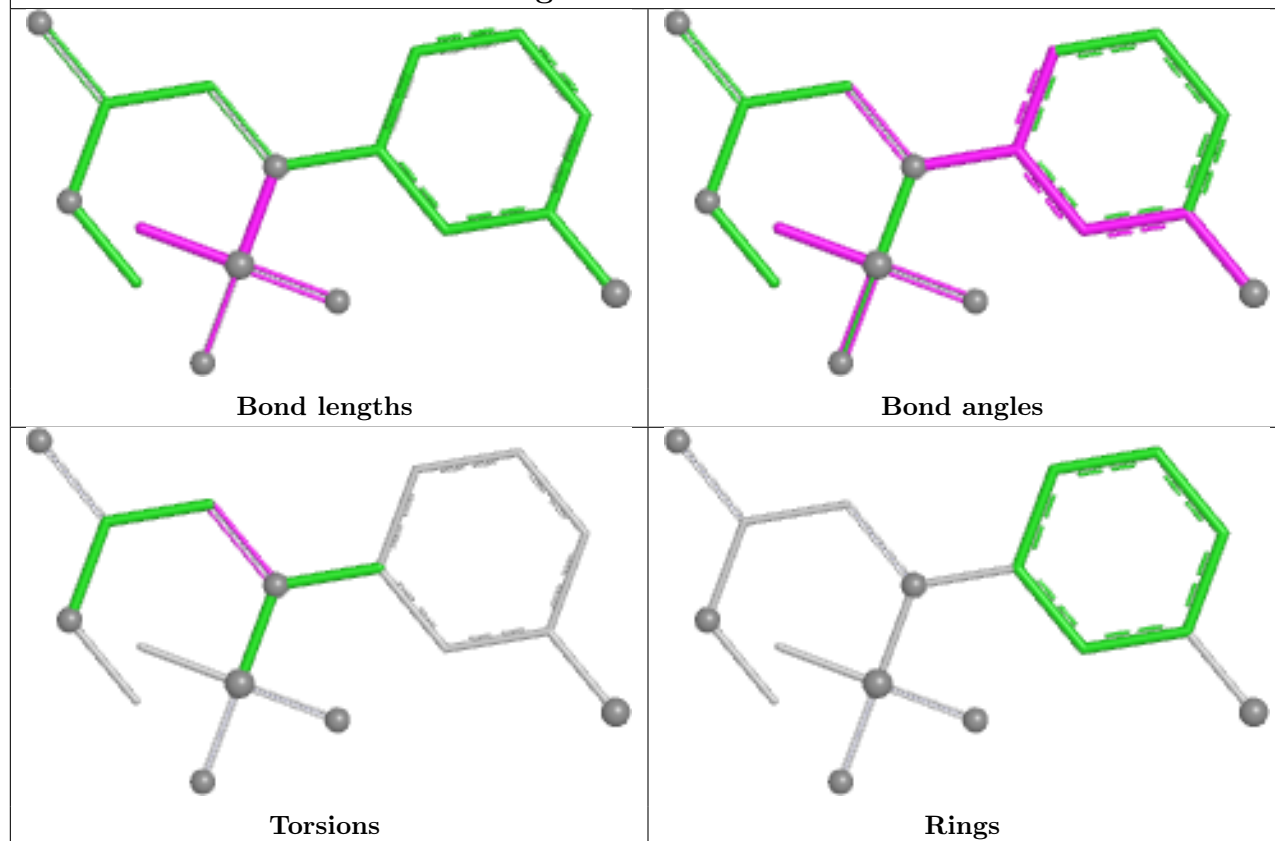
No monomer is involved in short contacts.

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.

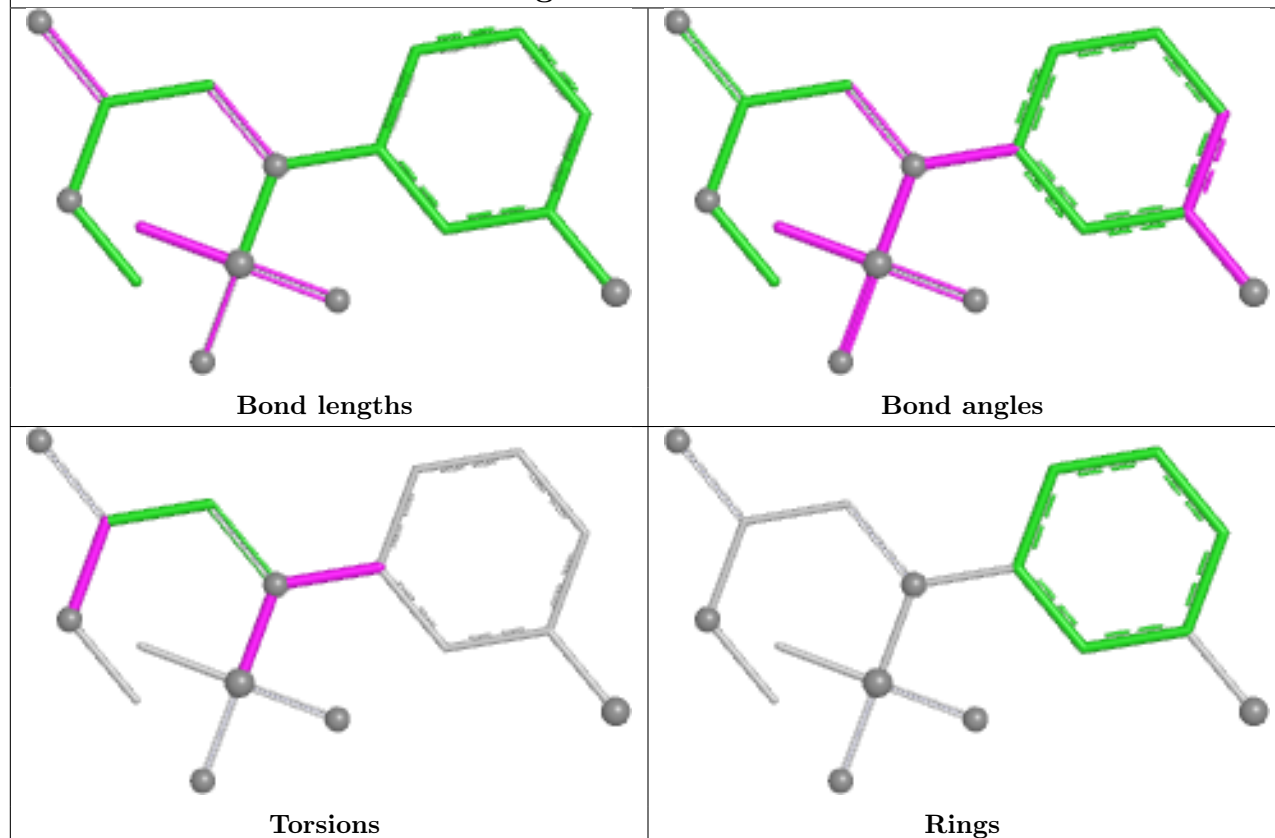
Ligand VBI B 401



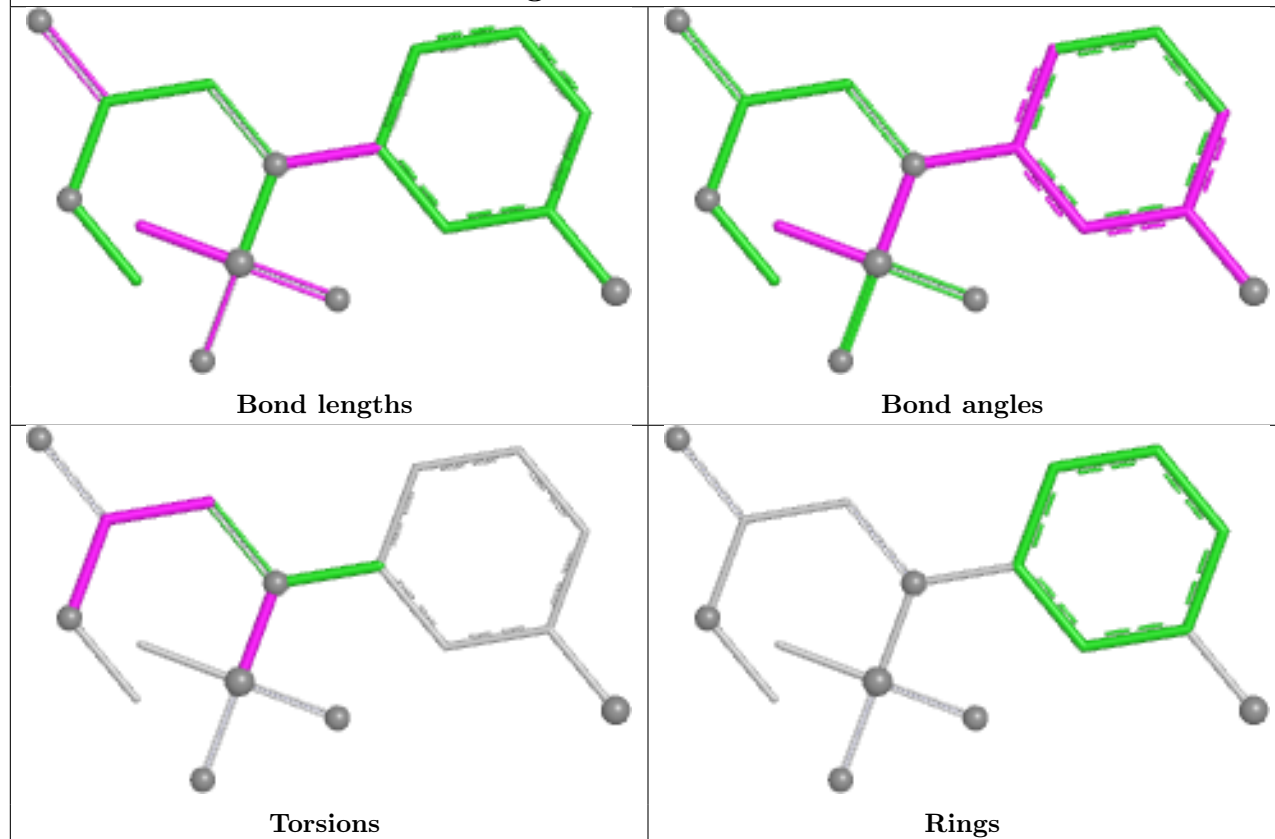
Ligand VBI B 403



Ligand VBI B 402



Ligand VBI A 2101



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	237/258 (91%)	1.21	46 (19%) 3 3	14, 41, 72, 88	12 (5%)
2	B	300/308 (97%)	1.49	77 (25%) 1 2	15, 44, 78, 104	8 (2%)
All	All	537/566 (94%)	1.37	123 (22%) 2 2	14, 42, 75, 104	20 (3%)

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	152	LEU	10.0
1	A	1979[A]	MET	8.3
2	B	170	SER	8.0
1	A	2069	VAL	7.7
2	B	7	THR	6.7
2	B	101	MET	5.9
2	B	41[A]	VAL	5.8
1	A	1833	GLY	5.7
1	A	2030	PRO	5.4
2	B	203[A]	TYR	5.3
2	B	108	ILE	5.2
2	B	279	TYR	5.0
2	B	316	LEU	5.0
2	B	317	LEU	4.9
2	B	21	SER	4.9
1	A	2027	LEU	4.8
2	B	172	PRO	4.8
2	B	173	ALA	4.8
1	A	2060	LEU	4.7
2	B	1	MET	4.6
2	B	84	ARG	4.5
1	A	2061	THR	4.4
2	B	275	LEU	4.4
2	B	29	PRO	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	1866	PHE	4.3
1	A	2037	TYR	4.2
2	B	55[A]	ARG	4.2
1	A	2064	GLY	4.2
2	B	282	ILE	4.1
1	A	2017[A]	THR	4.0
2	B	54	MET	4.0
1	A	2029	ASP	4.0
2	B	53	SER	3.9
1	A	1989	PHE	3.8
2	B	52	SER	3.8
2	B	111	ASP	3.8
2	B	106	PRO	3.8
2	B	110	GLU	3.8
1	A	2063	TYR	3.8
2	B	174	HIS	3.6
2	B	103	VAL	3.6
1	A	2068	ASN	3.5
2	B	258	LYS	3.5
2	B	91	ASN	3.5
1	A	2067	TYR	3.5
2	B	283	LEU	3.4
1	A	2049	ILE	3.3
2	B	276	PRO	3.3
1	A	1835	MET	3.3
2	B	87	ALA	3.3
1	A	1890	PHE	3.2
2	B	131	ARG	3.2
2	B	278	GLN	3.2
1	A	1834	ALA	3.2
1	A	2066	LYS	3.1
2	B	284	LEU	3.1
1	A	1840	TYR	3.1
2	B	96	PHE	3.1
1	A	1988	LEU	3.0
2	B	150	ASN	3.0
2	B	49	ALA	3.0
2	B	77	LEU	3.0
1	A	1962	ARG	3.0
2	B	107	LYS	3.0
1	A	1860	VAL	2.9
2	B	109	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	20	TYR	2.9
1	A	2028	SER	2.8
1	A	2031	THR	2.8
1	A	2058	LEU	2.8
2	B	100	GLN	2.8
1	A	2035	LYS	2.8
2	B	122[A]	GLN	2.8
2	B	68	TYR	2.8
2	B	234	PHE	2.7
2	B	147	VAL	2.7
2	B	38	ILE	2.7
1	A	1960[A]	GLU	2.7
1	A	2022	ALA	2.6
2	B	95	ASN	2.6
1	A	1862	VAL	2.6
2	B	51	ASN	2.6
2	B	17	ILE	2.6
2	B	171	ASP	2.6
1	A	2032	ILE	2.6
1	A	2033	THR	2.6
2	B	211[A]	GLN	2.6
1	A	1838	SER	2.5
2	B	277	GLU	2.5
2	B	213	ILE	2.5
1	A	1987	VAL	2.5
2	B	4	VAL	2.5
2	B	90	GLU	2.4
2	B	151	GLU	2.4
1	A	2065	ARG	2.4
1	A	2050	THR	2.4
2	B	88	LYS	2.4
1	A	2051	ILE	2.4
1	A	1931[A]	LYS	2.4
2	B	2	ASN	2.4
2	B	292	CYS	2.3
1	A	2053[A]	SER	2.3
2	B	235	GLY	2.3
2	B	62	CYS	2.3
2	B	315	GLU	2.3
1	A	2025	ILE	2.3
1	A	2024[A]	MET	2.3
1	A	1846	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	99	ARG	2.3
2	B	291	ILE	2.3
2	B	8	SER	2.3
1	A	1888	HIS	2.2
2	B	104	SER	2.2
2	B	124	ASP	2.2
2	B	89	PHE	2.2
2	B	233	PHE	2.2
2	B	254	ALA	2.1
2	B	313	TYR	2.1
1	A	1837	SER	2.1
2	B	212[A]	GLY	2.0
2	B	22	PHE	2.0
2	B	281	ASP	2.0
2	B	237	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

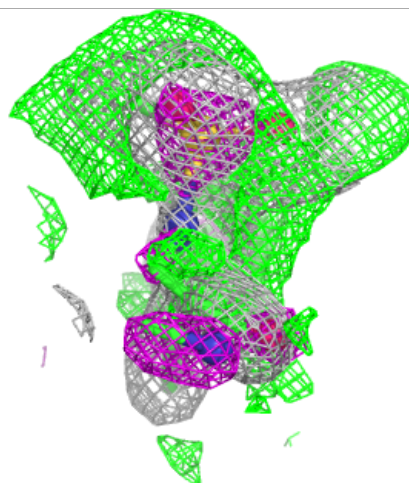
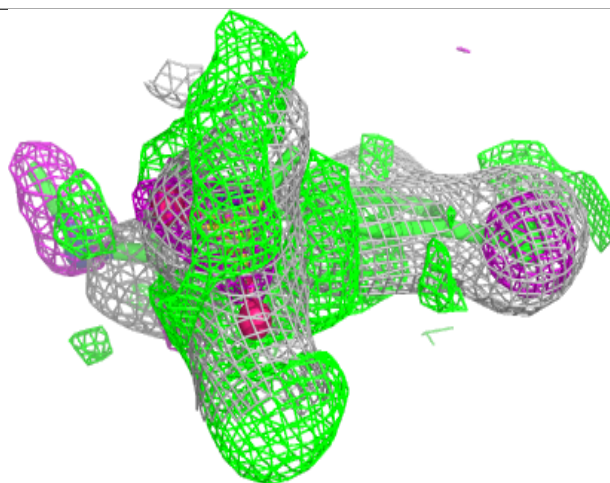
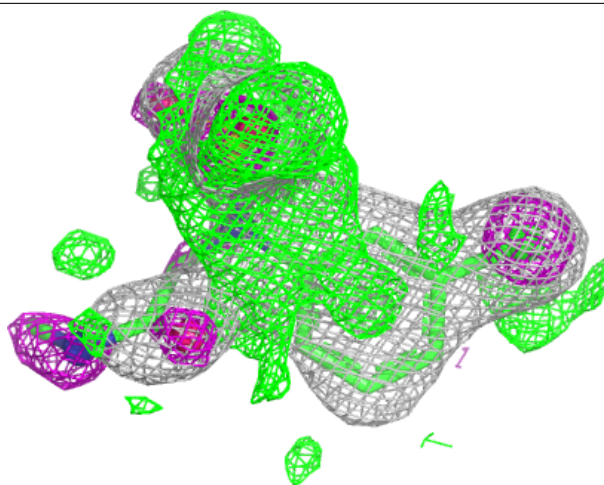
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	VBI	B	402	17/17	0.74	0.27	20,20,20,20	17
3	VBI	A	2101	17/17	0.76	0.25	20,20,20,20	17
3	VBI	B	401	17/17	0.79	0.22	20,20,20,20	17
3	VBI	B	403	17/17	0.85	0.23	20,20,20,20	17

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.

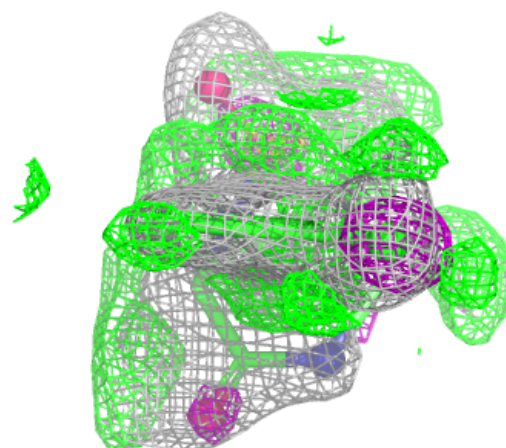
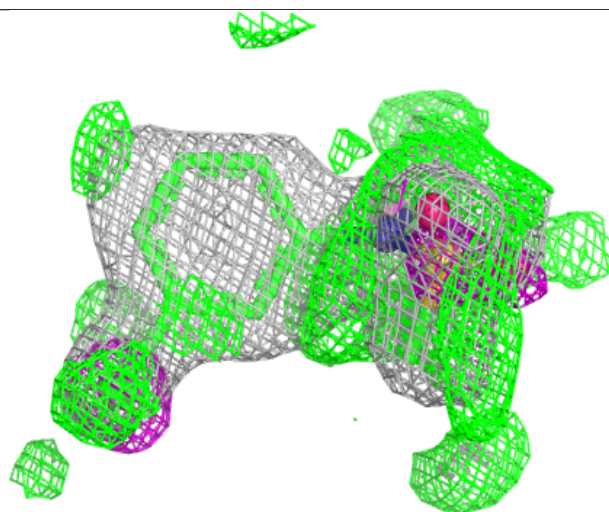
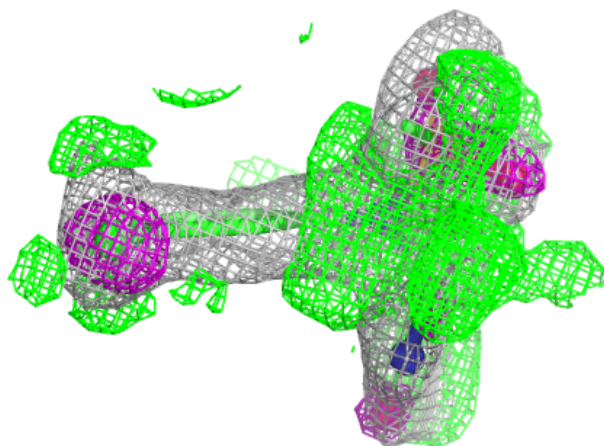
Electron density around VBI B 402:

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



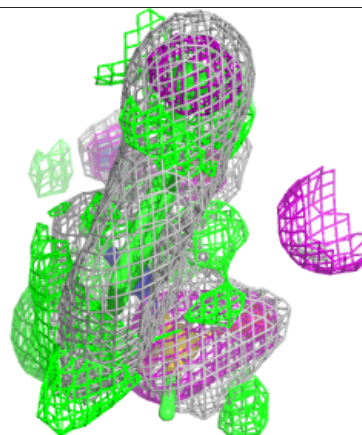
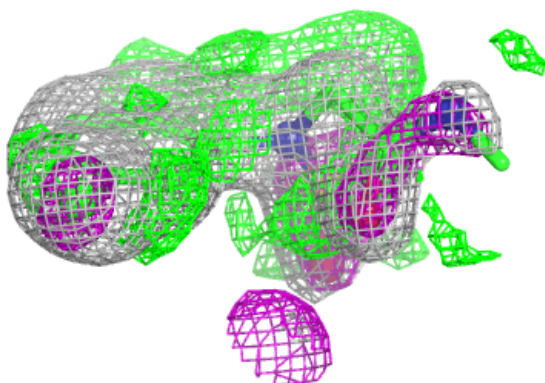
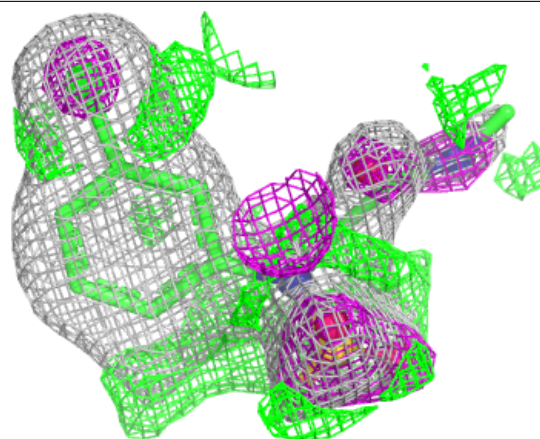
Electron density around VBI A 2101:

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



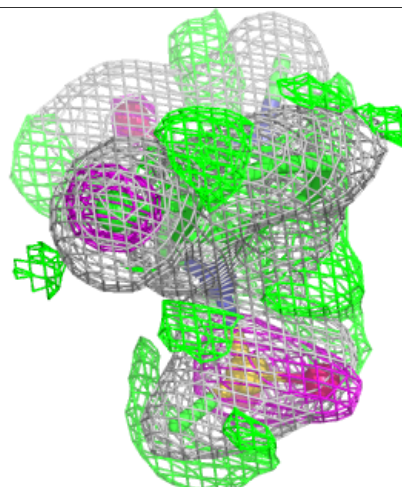
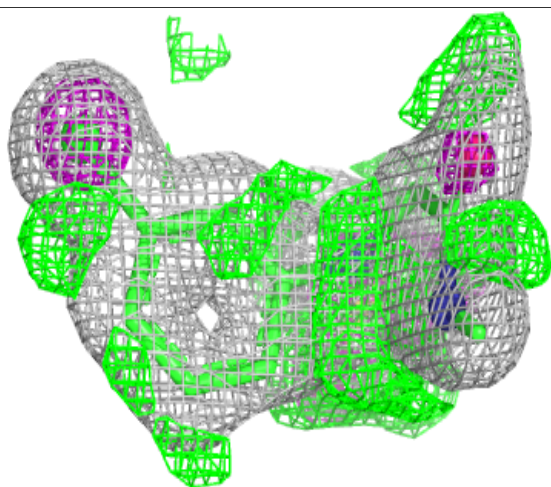
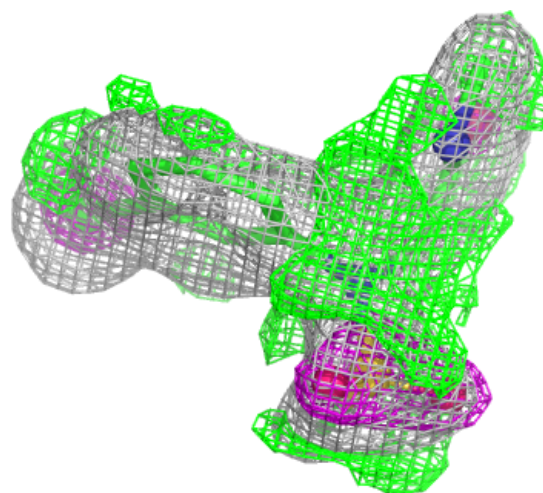
Electron density around VBI B 401:

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



Electron density around VBI B 403:

$2mF_o-DF_c$ (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.