



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:00 PM UTC

PDB ID : 4D3S / pdb_00004d3s
Title : Imine reductase from Nocardiosis halophila
Authors : Man, H.; Hart, S.; Turkenburg, J.P.; Grogan, G.
Deposited on : 2014-10-23
Resolution : 2.24 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

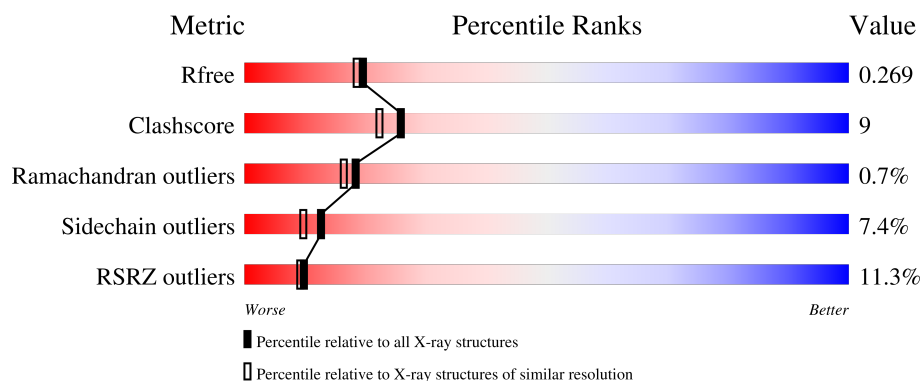
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	293	4% 80% 15% ..
1	B	293	6% 81% 14% ..
1	C	293	2% 81% 13% ...
1	D	293	3% 84% 12% ..
1	E	293	13% 75% 12% .. 9%

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Mol	Chain	Length	Quality of chain
1	F	293	4% 77% 16% • • •
1	G	293	32% 70% 17% • • 8%
1	H	293	22% 73% 18% • • 5%

2 Entry composition

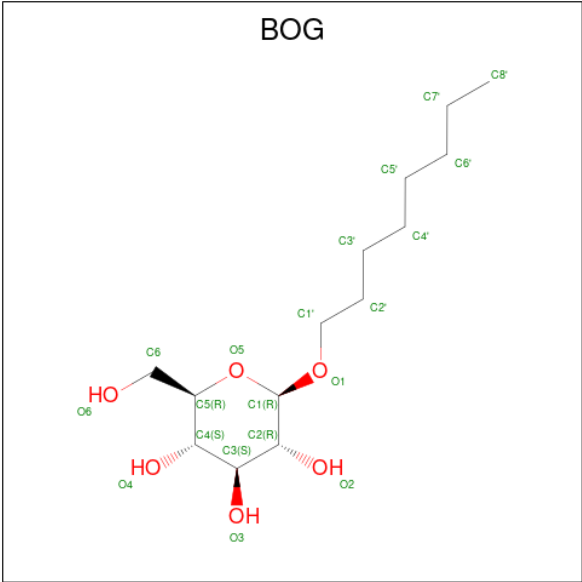
There are 3 unique types of molecules in this entry. The entry contains 16385 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 IMINE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2067	1303	352	402	10			
1	B	282	Total	C	N	O	S	0	0	0
			2013	1275	340	390	8			
1	C	283	Total	C	N	O	S	0	0	0
			2056	1296	346	404	10			
1	D	285	Total	C	N	O	S	0	0	0
			2047	1295	345	397	10			
1	E	267	Total	C	N	O	S	0	0	0
			1869	1188	310	361	10			
1	F	283	Total	C	N	O	S	0	0	0
			2036	1286	340	400	10			
1	G	269	Total	C	N	O	S	0	0	0
			1784	1119	312	343	10			
1	H	278	Total	C	N	O	S	0	0	0
			1966	1244	334	378	10			

- Molecule 2 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: C₁₄H₂₈O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			20	14	6		
2	H	1	Total	C	O	0	0
			20	14	6		

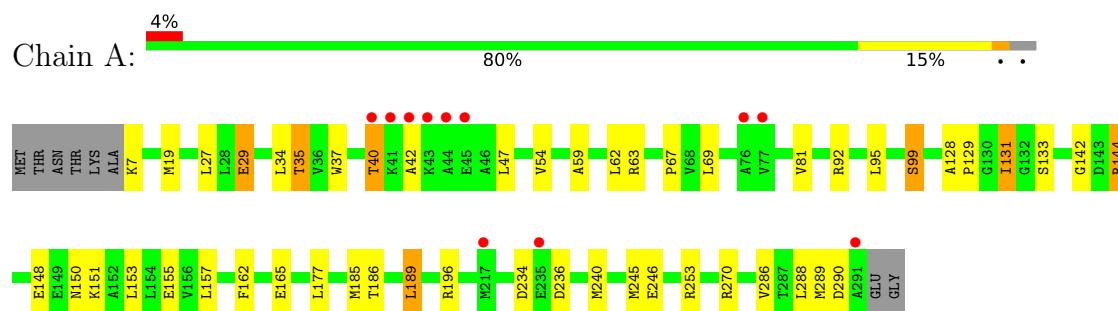
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	91	Total	O	0	0
			91	91		
3	B	73	Total	O	0	0
			73	73		
3	C	83	Total	O	0	0
			83	83		
3	D	97	Total	O	0	0
			97	97		
3	E	25	Total	O	0	0
			25	25		
3	F	69	Total	O	0	0
			69	69		
3	G	36	Total	O	0	0
			36	36		
3	H	33	Total	O	0	0
			33	33		

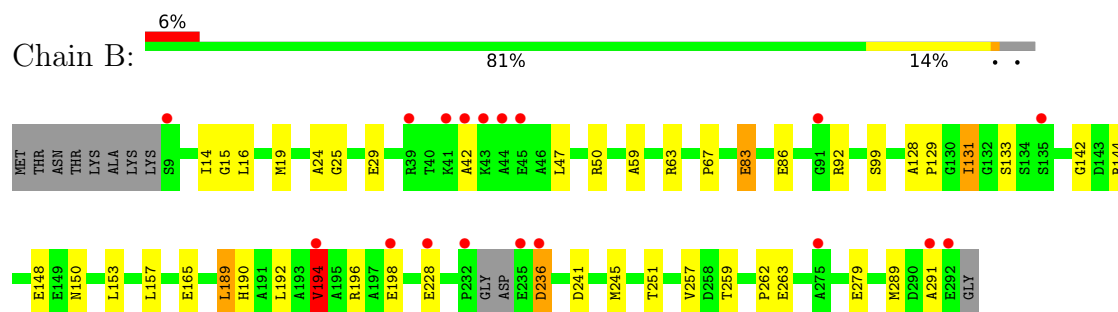
3 Residue-property plots

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.

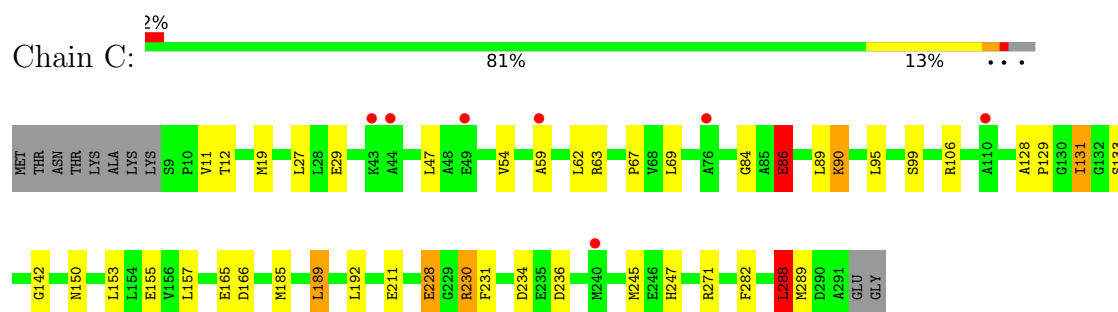
• Molecule 1: IMINE REDUCTASE



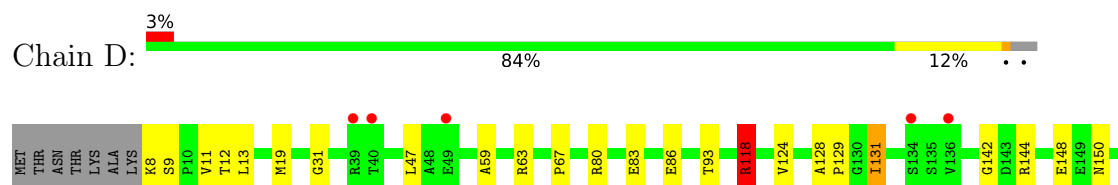
• Molecule 1: IMINE REDUCTASE

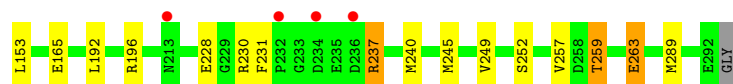


• Molecule 1: IMINE REDUCTASE

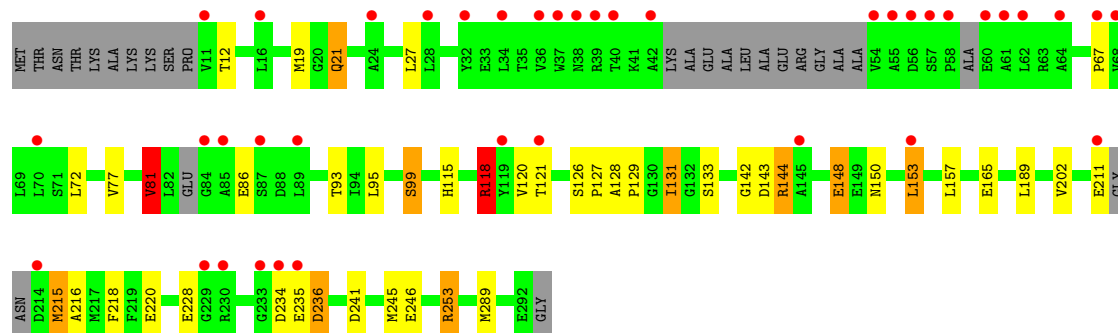
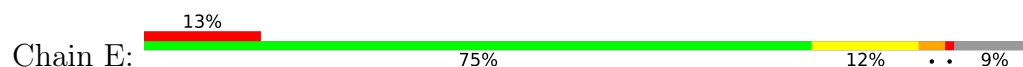


• Molecule 1: IMINE REDUCTASE

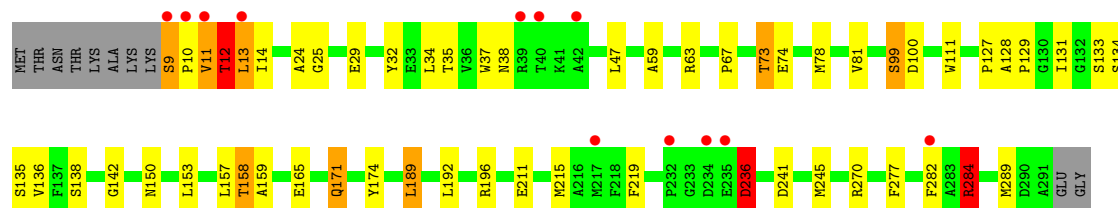
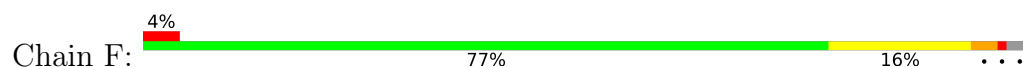




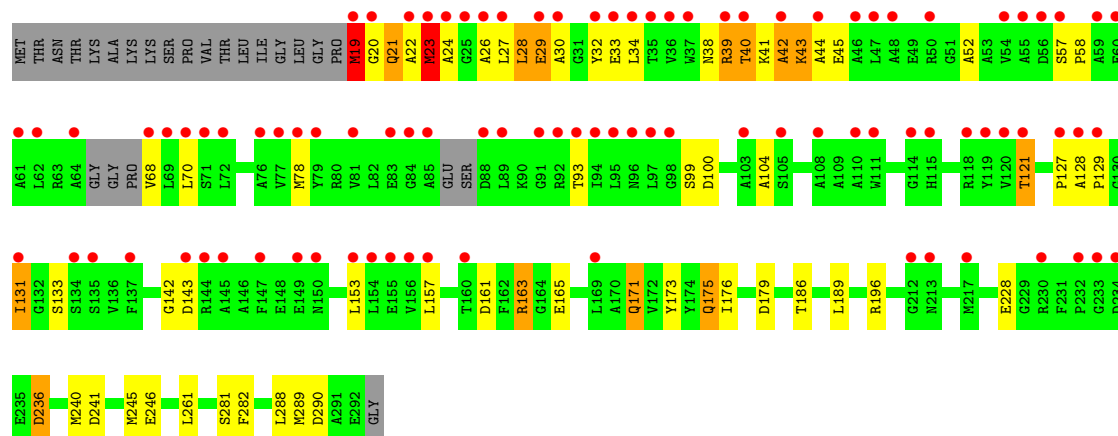
● Molecule 1: IMINE REDUCTASE



● Molecule 1: IMINE REDUCTASE

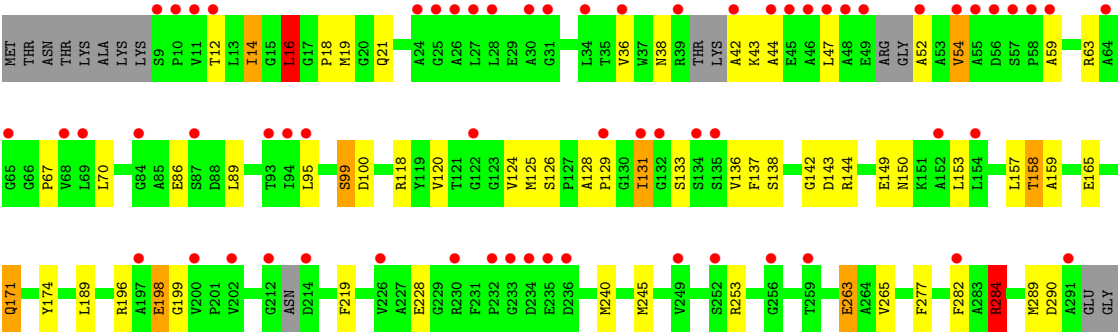


● Molecule 1: IMINE REDUCTASE



● Molecule 1: IMINE REDUCTASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	141.17Å 153.92Å 105.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.57 – 2.24 67.57 – 2.24	Depositor EDS
% Data completeness (in resolution range)	99.7 (67.57-2.24) 99.7 (67.57-2.24)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.230 , 0.266 0.235 , 0.269	Depositor DCC
R_{free} test set	5458 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	40.5	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16385	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 73.99 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6608e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BOG

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	0.93	1/2107 (0.0%)	0.96	5/2867 (0.2%)
1	B	0.89	0/2051	1.05	8/2796 (0.3%)
1	C	0.95	0/2096	0.97	8/2854 (0.3%)
1	D	0.91	0/2087	1.00	7/2843 (0.2%)
1	E	0.85	0/1902	0.98	9/2592 (0.3%)
1	F	1.08	3/2076 (0.1%)	1.26	17/2830 (0.6%)
1	G	0.94	1/1814 (0.1%)	1.03	7/2476 (0.3%)
1	H	0.88	0/2003	1.01	9/2728 (0.3%)
All	All	0.93	5/16136 (0.0%)	1.04	70/21986 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	F	0	2
All	All	0	3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	13	LEU	C-N	-24.01	1.00	1.33
1	G	121	THR	CA-CB	6.08	1.62	1.53
1	F	9	SER	C-N	5.56	1.40	1.33
1	A	40	THR	CA-C	5.56	1.60	1.52
1	F	10	PRO	N-CD	5.24	1.55	1.47

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	13	LEU	O-C-N	-19.83	100.92	123.33
1	F	13	LEU	CA-C-N	18.98	156.13	121.97
1	F	13	LEU	C-N-CA	18.98	156.13	121.97
1	B	14	ILE	CA-C-N	13.54	147.94	121.41
1	B	14	ILE	C-N-CA	13.54	147.94	121.41
1	D	118	ARG	CG-CD-NE	12.00	138.39	112.00
1	B	14	ILE	O-C-N	-11.88	110.43	123.26
1	F	12	THR	O-C-N	-11.51	108.60	123.21
1	F	11	VAL	CA-C-N	11.33	139.36	122.93
1	F	11	VAL	C-N-CA	11.33	139.36	122.93
1	H	284	ARG	CG-CD-NE	10.45	135.00	112.00
1	B	291	ALA	N-CA-C	9.66	121.50	110.97
1	F	133	SER	N-CA-C	9.62	122.74	110.33
1	G	133	SER	N-CA-C	9.36	121.92	110.41
1	H	133	SER	N-CA-C	8.15	120.43	110.41
1	F	14	ILE	CA-C-N	8.02	137.12	121.41
1	F	14	ILE	C-N-CA	8.02	137.12	121.41
1	E	234	ASP	N-CA-C	-7.74	99.02	110.48
1	E	228	GLU	N-CA-C	7.73	124.13	111.37
1	A	19	MET	CG-SD-CE	7.53	117.47	100.90
1	G	32	TYR	N-CA-C	7.48	121.67	109.85
1	E	21	GLN	N-CA-CB	7.42	120.77	110.01
1	H	228	GLU	N-CA-C	7.20	123.18	111.37
1	D	196	ARG	CG-CD-NE	7.15	127.74	112.00
1	G	23	MET	N-CA-CB	7.11	120.57	110.12
1	B	133	SER	N-CA-C	6.90	118.89	110.41
1	E	133	SER	N-CA-C	6.86	118.85	110.41
1	C	288	LEU	N-CA-CB	-6.84	99.25	110.40
1	G	228	GLU	N-CA-C	6.80	122.52	111.37
1	B	131	ILE	N-CA-C	6.63	117.38	110.62
1	C	133	SER	N-CA-C	6.63	118.56	110.41
1	F	9	SER	CA-C-N	-6.50	113.25	120.14
1	F	9	SER	C-N-CA	-6.50	113.25	120.14
1	F	284	ARG	CA-CB-CG	6.34	126.77	114.10
1	H	131	ILE	N-CA-C	6.29	117.04	110.62
1	H	284	ARG	NE-CZ-NH1	-6.29	115.21	121.50
1	D	11	VAL	CB-CA-C	-6.26	99.72	111.30
1	B	228	GLU	N-CA-C	6.20	121.53	111.37
1	C	230	ARG	N-CA-C	6.11	121.01	113.50
1	B	194	VAL	CB-CA-C	6.10	120.64	112.22
1	D	228	GLU	N-CA-C	6.04	121.27	111.37
1	A	131	ILE	N-CA-C	5.96	116.70	110.62
1	D	228	GLU	CB-CA-C	5.96	120.74	110.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	99	SER	CB-CA-C	5.95	119.56	109.50
1	D	118	ARG	CD-NE-CZ	5.92	132.69	124.40
1	C	11	VAL	CB-CA-C	-5.92	100.35	111.30
1	D	237	ARG	CA-CB-CG	5.83	125.77	114.10
1	C	131	ILE	N-CA-C	5.83	116.56	110.62
1	A	133	SER	N-CA-C	5.76	117.50	110.41
1	A	99	SER	CB-CA-C	5.71	118.94	109.53
1	E	118	ARG	CG-CD-NE	5.67	124.47	112.00
1	C	86	GLU	N-CA-CB	5.62	118.88	110.22
1	F	133	SER	CB-CA-C	-5.57	97.21	110.02
1	C	231	PHE	N-CA-C	5.47	121.91	109.81
1	H	16	LEU	CB-CA-C	-5.42	102.20	111.46
1	F	127	PRO	N-CA-C	-5.36	103.15	111.13
1	G	131	ILE	N-CA-C	5.34	116.07	110.62
1	F	236	ASP	N-CA-C	-5.34	100.20	108.42
1	E	131	ILE	N-CA-C	5.33	116.06	110.62
1	F	14	ILE	N-CA-C	5.30	120.37	109.34
1	G	121	THR	N-CA-CB	5.29	119.05	110.43
1	G	19	MET	N-CA-CB	5.28	119.47	110.50
1	H	199	GLY	CA-C-N	-5.20	118.85	122.59
1	H	199	GLY	C-N-CA	-5.20	118.85	122.59
1	C	155	GLU	CB-CG-CD	-5.20	103.76	112.60
1	F	11	VAL	CA-C-O	5.18	125.83	120.96
1	E	81	VAL	N-CA-C	5.17	120.10	109.34
1	A	92	ARG	CG-CD-NE	5.05	123.11	112.00
1	E	253	ARG	CG-CD-NE	5.01	123.02	112.00
1	H	43	LYS	N-CA-C	-5.00	106.46	113.37

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	15	GLY	Mainchain
1	F	12	THR	Mainchain
1	F	13	LEU	Mainchain

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	2067	0	2000	42	0
1	B	2013	0	1925	33	0
1	C	2056	0	1987	38	0
1	D	2047	0	1971	32	0
1	E	1869	0	1745	33	0
1	F	2036	0	1956	42	0
1	G	1784	0	1560	50	0
1	H	1966	0	1864	73	0
2	F	20	0	28	0	0
2	H	20	0	28	0	0
3	A	91	0	0	5	0
3	B	73	0	0	1	0
3	C	83	0	0	2	0
3	D	97	0	0	0	0
3	E	25	0	0	1	0
3	F	69	0	0	0	0
3	G	36	0	0	2	0
3	H	33	0	0	5	0
All	All	16385	0	15064	293	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 9.

All (293) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:185:MET:CE	1:F:215:MET:CE	2.24	1.15
1:H:16:LEU:HD11	1:H:38:ASN:HB2	1.21	1.11
1:C:185:MET:HE3	1:F:215:MET:HE1	1.26	1.11
1:C:185:MET:HE3	1:F:215:MET:CE	1.81	1.10
1:E:77:VAL:O	1:E:81:VAL:HG12	1.54	1.07
1:E:215:MET:CE	1:E:215:MET:HA	1.90	1.02
1:H:126:SER:HB2	1:H:136:VAL:HG21	1.00	0.98
1:H:126:SER:HB2	1:H:136:VAL:CG2	1.94	0.98
1:H:16:LEU:HD11	1:H:38:ASN:CB	1.97	0.95
1:H:126:SER:CB	1:H:136:VAL:HG21	1.96	0.94
1:C:185:MET:CE	1:F:215:MET:HE3	1.96	0.93
1:H:136:VAL:HG13	1:H:158:THR:OG1	1.67	0.93
1:G:100:ASP:H	1:G:171:GLN:HE21	1.18	0.92
1:H:16:LEU:HD12	1:H:38:ASN:ND2	1.89	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:MET:HE1	1:H:18:PRO:CD	2.04	0.87
1:E:215:MET:HA	1:E:215:MET:HE2	1.55	0.87
3:A:2043:HOH:O	1:H:198:GLU:HG2	1.74	0.87
1:G:186:THR:HG22	1:G:261:LEU:HD22	1.57	0.86
1:C:185:MET:HE1	1:F:215:MET:HE3	1.56	0.85
1:H:158:THR:HG23	1:H:159:ALA:O	1.77	0.85
1:G:28:LEU:O	1:G:30:ALA:N	2.09	0.85
1:A:290:ASP:OD2	1:H:196:ARG:NH1	2.11	0.82
1:H:263:GLU:HG3	3:H:2030:HOH:O	1.81	0.79
1:E:215:MET:HA	1:E:215:MET:HE3	1.66	0.76
1:F:158:THR:HG23	1:F:159:ALA:O	1.84	0.76
1:G:99:SER:HA	1:G:171:GLN:NE2	2.02	0.75
1:A:189:LEU:HB3	1:H:289:MET:HE1	1.69	0.73
1:H:284:ARG:CG	1:H:284:ARG:HH11	2.02	0.72
1:D:252:SER:O	1:D:257:VAL:HG12	1.90	0.72
1:F:11:VAL:O	1:F:34:LEU:HA	1.89	0.72
1:D:128:ALA:HA	1:D:131:ILE:HD12	1.70	0.71
1:G:99:SER:HA	1:G:171:GLN:HE22	1.55	0.71
1:H:16:LEU:CD1	1:H:38:ASN:CG	2.63	0.71
1:C:289:MET:HE1	1:F:189:LEU:HB3	1.74	0.70
1:D:93:THR:OG1	1:D:118:ARG:NH1	2.25	0.69
1:B:144:ARG:NH1	1:B:148:GLU:OE2	2.25	0.69
1:F:100:ASP:O	1:F:171:GLN:HG3	1.92	0.69
1:C:106:ARG:NH1	1:C:166:ASP:OD2	2.26	0.68
1:G:21:GLN:O	1:G:24:ALA:N	2.26	0.68
1:D:230:ARG:HG3	1:D:231:PHE:H	1.59	0.68
1:D:144:ARG:NH1	1:D:148:GLU:OE2	2.26	0.67
1:H:100:ASP:O	1:H:171:GLN:HG3	1.95	0.66
1:H:16:LEU:CD1	1:H:38:ASN:ND2	2.57	0.66
1:G:19:MET:CE	1:G:128:ALA:HA	2.26	0.66
1:E:93:THR:OG1	1:E:118:ARG:NH1	2.28	0.65
1:H:14:ILE:HG23	1:H:70:LEU:HD23	1.78	0.65
1:B:189:LEU:HB3	1:G:289:MET:HE1	1.78	0.65
1:B:190:HIS:HD2	1:G:179:ASP:OD2	1.80	0.65
1:G:38:ASN:O	1:G:40:THR:N	2.30	0.64
1:H:284:ARG:HH11	1:H:284:ARG:HG2	1.62	0.64
1:H:16:LEU:CD1	1:H:38:ASN:HB2	2.14	0.64
1:C:189:LEU:HD22	1:F:282:PHE:CZ	2.33	0.64
1:G:19:MET:HE1	1:G:128:ALA:N	2.13	0.64
1:F:136:VAL:HG23	1:F:158:THR:OG1	1.96	0.64
1:C:185:MET:CE	1:F:215:MET:HE1	2.04	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:277:PHE:CE1	1:H:284:ARG:NH1	2.66	0.64
1:G:186:THR:HG22	1:G:261:LEU:CD2	2.27	0.63
1:H:95:LEU:HD23	1:H:120:VAL:HG22	1.79	0.63
1:G:43:LYS:C	1:G:45:GLU:H	2.07	0.62
1:H:158:THR:HG22	3:H:2021:HOH:O	2.00	0.62
1:A:240:MET:HE1	1:H:18:PRO:CG	2.30	0.61
1:G:19:MET:HE2	1:G:128:ALA:HA	1.82	0.61
1:E:215:MET:CE	1:E:218:PHE:CD2	2.83	0.61
1:E:95:LEU:HD23	1:E:120:VAL:HG22	1.81	0.61
1:A:35:THR:CG2	3:A:2003:HOH:O	2.49	0.61
1:C:84:GLY:N	1:C:86:GLU:OE2	2.33	0.60
1:B:194:VAL:CG2	1:G:176:ILE:HG12	2.31	0.60
1:F:138:SER:OG	1:F:158:THR:HG21	2.01	0.60
1:A:189:LEU:HD22	1:H:282:PHE:CZ	2.38	0.59
1:H:16:LEU:CD1	1:H:38:ASN:CB	2.77	0.59
1:C:131:ILE:O	1:C:157:LEU:O	2.21	0.59
1:D:237:ARG:HG2	1:D:240:MET:HE3	1.84	0.59
1:G:100:ASP:H	1:G:171:GLN:NE2	1.93	0.59
1:B:259:THR:HG23	1:B:263:GLU:OE2	2.03	0.59
1:C:86:GLU:HA	1:C:89:LEU:HD13	1.85	0.59
1:A:289:MET:HE1	1:H:189:LEU:HB3	1.84	0.59
1:B:190:HIS:CD2	1:G:179:ASP:OD2	2.57	0.58
1:C:189:LEU:HB3	1:F:289:MET:HE1	1.85	0.58
1:G:23:MET:O	1:G:26:ALA:N	2.36	0.58
1:C:90:LYS:HD3	1:C:90:LYS:N	2.17	0.58
1:C:106:ARG:NH1	1:C:166:ASP:OD1	2.36	0.58
1:B:131:ILE:O	1:B:157:LEU:O	2.22	0.58
1:E:19:MET:HE1	1:E:127:PRO:HA	1.86	0.58
1:A:37:TRP:CZ3	1:A:81:VAL:CG2	2.87	0.58
1:A:40:THR:OG1	1:A:42:ALA:N	2.36	0.58
1:G:131:ILE:O	1:G:157:LEU:O	2.22	0.58
1:F:25:GLY:O	1:F:29:GLU:HG3	2.03	0.57
1:E:131:ILE:O	1:E:157:LEU:O	2.22	0.57
1:D:230:ARG:CG	1:D:231:PHE:H	2.18	0.57
1:E:215:MET:CE	1:E:215:MET:CA	2.74	0.57
1:A:131:ILE:O	1:A:157:LEU:O	2.23	0.56
1:H:42:ALA:C	1:H:44:ALA:H	2.11	0.56
1:H:86:GLU:HA	1:H:89:LEU:HD13	1.86	0.56
1:C:90:LYS:HD3	1:C:90:LYS:O	2.05	0.56
1:G:39:ARG:O	1:G:41:LYS:N	2.39	0.56
1:A:40:THR:OG1	1:A:42:ALA:HB2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:GLY:O	1:B:29:GLU:HG3	2.05	0.56
1:F:131:ILE:O	1:F:157:LEU:O	2.24	0.56
1:A:240:MET:HE1	1:H:18:PRO:HD3	1.85	0.55
1:B:19:MET:HE1	3:B:2038:HOH:O	2.05	0.55
1:F:73:THR:HG22	1:F:74:GLU:HG2	1.88	0.55
1:G:27:LEU:O	1:G:28:LEU:O	2.25	0.55
1:C:247:HIS:HD2	3:C:2069:HOH:O	1.88	0.55
1:F:158:THR:CG2	1:F:159:ALA:O	2.54	0.55
1:D:259:THR:HG23	1:D:263:GLU:OE1	2.06	0.55
1:H:131:ILE:O	1:H:157:LEU:O	2.25	0.55
1:E:86:GLU:CB	1:E:115:HIS:CD2	2.89	0.55
1:E:215:MET:CE	1:E:218:PHE:HD2	2.20	0.55
1:D:144:ARG:HH12	1:D:148:GLU:CD	2.14	0.55
1:D:237:ARG:CG	1:D:240:MET:HE3	2.37	0.55
1:D:252:SER:HB3	1:D:257:VAL:HG13	1.87	0.55
1:G:70:LEU:CD1	1:G:78:MET:HE3	2.37	0.54
1:H:131:ILE:CB	3:H:2020:HOH:O	2.55	0.54
1:A:67:PRO:HG3	1:A:150:ASN:OD1	2.07	0.54
1:A:37:TRP:CZ3	1:A:81:VAL:HG23	2.41	0.54
1:C:19:MET:HE1	3:C:2034:HOH:O	2.08	0.54
1:C:271:ARG:NH2	1:C:288:LEU:HD23	2.22	0.54
1:H:138:SER:OG	1:H:158:THR:HG21	2.07	0.54
1:C:27:LEU:CD2	1:C:153:LEU:HD11	2.38	0.53
1:H:16:LEU:HD11	1:H:38:ASN:CG	2.30	0.53
1:A:185:MET:HE1	1:H:219:PHE:CE1	2.43	0.53
1:D:237:ARG:HG2	1:D:240:MET:CE	2.38	0.53
1:G:19:MET:CE	1:G:128:ALA:CA	2.87	0.53
1:B:67:PRO:HG3	1:B:150:ASN:OD1	2.08	0.53
1:A:40:THR:OG1	1:A:42:ALA:CB	2.56	0.53
1:E:67:PRO:HG3	1:E:150:ASN:OD1	2.08	0.53
1:C:106:ARG:NH1	1:C:166:ASP:CG	2.67	0.53
1:D:67:PRO:HG3	1:D:150:ASN:OD1	2.08	0.53
1:H:67:PRO:HG3	1:H:150:ASN:OD1	2.09	0.53
1:H:99:SER:HA	3:H:2017:HOH:O	2.08	0.53
1:C:282:PHE:CZ	1:F:189:LEU:HD22	2.44	0.52
1:A:144:ARG:O	1:A:148:GLU:OE1	2.26	0.52
1:H:253:ARG:NH2	1:H:263:GLU:OE2	2.43	0.52
1:C:90:LYS:O	1:C:90:LYS:HE2	2.09	0.52
1:D:8:LYS:HG2	1:D:31:GLY:O	2.10	0.52
1:F:67:PRO:HG3	1:F:150:ASN:OD1	2.08	0.52
1:E:215:MET:HE1	1:E:218:PHE:CD2	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ARG:HH12	1:B:148:GLU:CD	2.16	0.52
1:C:67:PRO:HG3	1:C:150:ASN:OD1	2.09	0.52
1:C:289:MET:CE	1:F:189:LEU:HB3	2.39	0.52
1:G:281:SER:OG	1:G:282:PHE:N	2.42	0.52
1:H:158:THR:CG2	1:H:159:ALA:O	2.56	0.52
1:C:185:MET:HE1	1:F:219:PHE:CE1	2.44	0.52
1:G:21:GLN:O	1:G:23:MET:N	2.42	0.52
1:D:8:LYS:CG	1:D:31:GLY:O	2.57	0.52
1:H:277:PHE:CZ	1:H:284:ARG:NH1	2.78	0.52
1:E:202:VAL:HG23	3:E:2014:HOH:O	2.10	0.51
1:G:171:GLN:HA	1:G:171:GLN:OE1	2.09	0.51
1:A:240:MET:CE	1:H:18:PRO:CD	2.83	0.51
1:F:192:LEU:O	1:F:196:ARG:HG3	2.11	0.51
1:B:128:ALA:HB3	1:B:129:PRO:HD3	1.93	0.51
1:E:27:LEU:CD2	1:E:153:LEU:HD11	2.39	0.51
1:A:151:LYS:O	1:A:155:GLU:HG3	2.11	0.51
1:B:189:LEU:HD22	1:G:282:PHE:CZ	2.46	0.51
1:C:27:LEU:HD22	1:C:153:LEU:HD11	1.92	0.51
1:A:253:ARG:HD3	1:B:92:ARG:NH1	2.26	0.50
1:F:37:TRP:CZ3	1:F:81:VAL:CG2	2.94	0.50
1:A:290:ASP:CG	1:H:196:ARG:HH11	2.19	0.50
1:C:27:LEU:CD2	1:C:153:LEU:CD1	2.89	0.50
1:C:228:GLU:OE2	1:C:230:ARG:NH1	2.44	0.50
1:A:240:MET:HE1	1:H:18:PRO:HD2	1.93	0.50
1:D:252:SER:HB3	1:D:257:VAL:CG1	2.41	0.50
1:E:27:LEU:CD2	1:E:153:LEU:CD1	2.90	0.50
1:F:24:ALA:HB1	1:F:47:LEU:HD22	1.94	0.50
1:A:196:ARG:NH1	1:H:290:ASP:OD2	2.45	0.50
1:G:163:ARG:HD3	1:G:173:TYR:CZ	2.47	0.50
1:H:95:LEU:CD2	1:H:120:VAL:CG2	2.90	0.50
1:D:144:ARG:NH1	1:D:148:GLU:CG	2.75	0.49
1:G:161:ASP:OD1	1:G:163:ARG:NH2	2.37	0.49
1:H:136:VAL:HG22	1:H:137:PHE:N	2.28	0.49
1:B:236:ASP:OD1	1:B:241:ASP:OD2	2.30	0.49
1:D:144:ARG:NH1	1:D:148:GLU:HG2	2.28	0.49
1:E:216:ALA:O	1:E:220:GLU:HG3	2.13	0.49
1:A:289:MET:HE1	1:H:189:LEU:CB	2.43	0.49
1:H:16:LEU:HD21	1:H:36:VAL:CG1	2.43	0.49
1:G:41:LYS:O	1:G:42:ALA:HB3	2.12	0.49
1:A:128:ALA:HB3	1:A:129:PRO:HD3	1.94	0.49
1:C:228:GLU:O	1:C:228:GLU:HG3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:27:LEU:HD22	1:E:153:LEU:HD11	1.95	0.49
1:E:118:ARG:HG3	1:E:118:ARG:HH11	1.77	0.49
1:B:289:MET:HE1	1:G:189:LEU:HB3	1.95	0.48
1:D:128:ALA:HB3	1:D:129:PRO:HD3	1.94	0.48
1:F:236:ASP:OD1	1:F:241:ASP:OD2	2.30	0.48
1:C:69:LEU:HD23	1:C:95:LEU:HB2	1.96	0.48
1:D:257:VAL:HG23	1:E:289:MET:O	2.13	0.48
1:D:289:MET:HE1	1:E:189:LEU:HB3	1.96	0.48
1:F:11:VAL:HG12	1:F:32:TYR:HB3	1.95	0.48
1:A:29:GLU:HG3	3:A:2010:HOH:O	2.14	0.48
1:A:289:MET:CE	1:H:189:LEU:HB3	2.43	0.48
1:G:43:LYS:C	1:G:45:GLU:N	2.70	0.48
1:A:69:LEU:HD23	1:A:95:LEU:HB2	1.95	0.47
1:C:90:LYS:O	1:C:90:LYS:CD	2.62	0.47
1:F:128:ALA:HB3	1:F:129:PRO:HD3	1.96	0.47
1:C:128:ALA:HB3	1:C:129:PRO:HD3	1.96	0.47
1:D:8:LYS:CG	1:D:9:SER:H	2.27	0.47
1:E:128:ALA:HB3	1:E:129:PRO:HD3	1.96	0.47
1:G:28:LEU:O	1:G:29:GLU:C	2.55	0.47
1:B:144:ARG:NH1	1:B:148:GLU:HG2	2.30	0.47
1:F:37:TRP:CZ3	1:F:81:VAL:HG23	2.50	0.47
1:E:95:LEU:CD2	1:E:120:VAL:CG2	2.93	0.47
1:C:185:MET:HE2	1:F:215:MET:CE	2.36	0.47
1:H:42:ALA:C	1:H:44:ALA:N	2.73	0.47
1:E:236:ASP:OD2	1:E:241:ASP:OD2	2.33	0.46
1:E:19:MET:HE1	1:E:126:SER:O	2.14	0.46
1:F:100:ASP:O	1:F:171:GLN:CG	2.60	0.46
1:G:236:ASP:OD2	1:G:241:ASP:OD2	2.33	0.46
1:E:120:VAL:HG12	1:E:143:ASP:N	2.29	0.46
1:H:120:VAL:HG12	1:H:143:ASP:N	2.30	0.46
1:A:240:MET:CE	1:H:18:PRO:HD2	2.46	0.46
1:B:189:LEU:HB3	1:G:289:MET:CE	2.46	0.46
1:H:128:ALA:HB3	1:H:129:PRO:HD3	1.97	0.46
1:A:40:THR:HG1	1:A:42:ALA:N	2.13	0.46
1:B:194:VAL:HG12	1:B:257:VAL:HG21	1.98	0.46
1:A:153:LEU:HD23	1:A:153:LEU:C	2.41	0.46
1:B:83:GLU:HG3	1:D:80:ARG:NE	2.30	0.46
1:G:128:ALA:HB3	1:G:129:PRO:HD3	1.97	0.46
1:H:47:LEU:O	1:H:52:ALA:HB1	2.15	0.46
1:B:24:ALA:HB1	1:B:47:LEU:HD22	1.97	0.46
1:A:286:VAL:HB	1:H:189:LEU:HD22	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:100:ASP:OD2	1:G:104:ALA:CB	2.64	0.45
1:B:251:THR:OG1	1:G:175:GLN:HG2	2.15	0.45
1:G:41:LYS:O	1:G:42:ALA:CB	2.64	0.45
1:D:142:GLY:O	1:D:165:GLU:HA	2.17	0.45
1:A:37:TRP:CH2	1:A:81:VAL:HG23	2.52	0.45
1:B:142:GLY:O	1:B:165:GLU:HA	2.17	0.45
1:B:144:ARG:NH1	1:B:148:GLU:CG	2.79	0.45
1:H:54:VAL:O	1:H:54:VAL:CG2	2.65	0.45
1:B:196:ARG:NH1	1:G:290:ASP:OD2	2.48	0.44
1:F:99:SER:HB2	1:F:174:TYR:CE2	2.52	0.44
1:G:19:MET:HE1	1:G:127:PRO:C	2.42	0.44
1:H:100:ASP:O	1:H:171:GLN:CG	2.62	0.44
1:C:142:GLY:O	1:C:165:GLU:HA	2.17	0.44
1:A:189:LEU:HB3	1:H:289:MET:CE	2.41	0.44
1:B:194:VAL:HG22	1:G:176:ILE:HG12	1.97	0.44
1:F:37:TRP:CH2	1:F:81:VAL:HG23	2.53	0.44
1:D:230:ARG:HG3	1:D:231:PHE:N	2.28	0.44
1:H:16:LEU:HD12	1:H:38:ASN:CG	2.33	0.44
1:E:144:ARG:NE	1:E:148:GLU:OE1	2.50	0.44
1:G:153:LEU:C	1:G:153:LEU:HD23	2.43	0.44
1:A:37:TRP:CZ3	1:A:81:VAL:HG21	2.52	0.44
1:H:118:ARG:HH22	1:H:149:GLU:CD	2.26	0.43
1:H:142:GLY:O	1:H:165:GLU:HA	2.17	0.43
1:G:196:ARG:HA	3:G:2012:HOH:O	2.17	0.43
1:H:153:LEU:HD23	1:H:153:LEU:C	2.43	0.43
1:F:153:LEU:C	1:F:153:LEU:HD23	2.43	0.43
1:H:59:ALA:O	1:H:63:ARG:HG3	2.18	0.43
1:A:35:THR:HG22	3:A:2003:HOH:O	2.16	0.43
1:B:59:ALA:O	1:B:63:ARG:HG3	2.19	0.43
1:F:142:GLY:O	1:F:165:GLU:HA	2.19	0.43
1:A:186:THR:HG21	1:H:265:VAL:HG13	2.00	0.42
1:C:59:ALA:O	1:C:63:ARG:HG3	2.20	0.42
1:B:153:LEU:C	1:B:153:LEU:HD23	2.44	0.42
1:G:38:ASN:O	1:G:39:ARG:C	2.61	0.42
1:H:16:LEU:HD22	1:H:44:ALA:HB2	2.00	0.42
1:F:277:PHE:CE1	1:F:284:ARG:NH1	2.87	0.42
1:G:142:GLY:O	1:G:165:GLU:HA	2.19	0.42
1:D:59:ALA:O	1:D:63:ARG:HG3	2.20	0.42
1:B:83:GLU:HG3	1:D:80:ARG:CD	2.50	0.42
1:B:83:GLU:O	1:B:86:GLU:OE1	2.38	0.42
1:D:128:ALA:HA	1:D:131:ILE:CD1	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:95:LEU:HD23	1:E:120:VAL:CG2	2.49	0.42
1:A:142:GLY:O	1:A:165:GLU:HA	2.20	0.42
1:B:50:ARG:HD3	1:F:134:SER:HB3	2.02	0.42
1:D:153:LEU:C	1:D:153:LEU:HD23	2.45	0.42
1:F:59:ALA:O	1:F:63:ARG:HG3	2.19	0.42
1:H:95:LEU:CD2	1:H:120:VAL:HG22	2.48	0.42
1:D:19:MET:CE	1:D:124:VAL:HG11	2.50	0.41
1:E:95:LEU:CD2	1:E:120:VAL:HG22	2.49	0.41
1:G:68:VAL:CB	1:G:93:THR:O	2.68	0.41
1:H:99:SER:HB2	1:H:174:TYR:CE2	2.55	0.41
1:B:190:HIS:NE2	1:B:262:PRO:HG3	2.35	0.41
1:F:12:THR:HA	1:F:35:THR:O	2.20	0.41
1:G:20:GLY:O	1:G:21:GLN:C	2.63	0.41
1:A:144:ARG:HD2	1:A:162:PHE:CE2	2.55	0.41
1:E:142:GLY:O	1:E:165:GLU:HA	2.20	0.41
1:A:29:GLU:CG	3:A:2010:HOH:O	2.67	0.41
1:C:185:MET:CE	1:F:215:MET:HE2	2.40	0.41
1:H:125:MET:O	1:H:136:VAL:HG23	2.21	0.41
1:B:190:HIS:O	1:B:194:VAL:HG13	2.20	0.41
1:C:189:LEU:HB3	1:F:289:MET:CE	2.50	0.41
1:E:19:MET:HE1	1:E:127:PRO:CA	2.51	0.41
1:F:78:MET:HE3	1:F:111:TRP:CE3	2.56	0.41
1:H:95:LEU:HD22	1:H:120:VAL:CG2	2.51	0.41
1:D:83:GLU:O	1:D:86:GLU:OE1	2.39	0.41
1:E:77:VAL:O	1:E:81:VAL:CG1	2.46	0.41
1:G:143:ASP:CB	3:G:2008:HOH:O	2.69	0.41
1:H:240:MET:HG3	3:H:2028:HOH:O	2.20	0.40
1:A:240:MET:HE1	1:H:18:PRO:HG2	2.00	0.40
1:H:136:VAL:CG1	1:H:158:THR:OG1	2.53	0.40
1:A:59:ALA:O	1:A:63:ARG:HG3	2.22	0.40
1:G:21:GLN:H	1:G:21:GLN:HG2	1.62	0.40
1:D:257:VAL:HG12	1:D:257:VAL:H	1.65	0.40
1:H:19:MET:CE	1:H:124:VAL:HG11	2.50	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	283/293 (97%)	279 (99%)	4 (1%)	0	100	100
1	B	278/293 (95%)	272 (98%)	5 (2%)	1 (0%)	30	30
1	C	281/293 (96%)	278 (99%)	3 (1%)	0	100	100
1	D	283/293 (97%)	280 (99%)	3 (1%)	0	100	100
1	E	257/293 (88%)	250 (97%)	6 (2%)	1 (0%)	30	30
1	F	281/293 (96%)	277 (99%)	4 (1%)	0	100	100
1	G	263/293 (90%)	240 (91%)	10 (4%)	13 (5%)	1	0
1	H	270/293 (92%)	267 (99%)	3 (1%)	0	100	100
All	All	2196/2344 (94%)	2143 (98%)	38 (2%)	15 (1%)	18	16

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	28	LEU
1	G	29	GLU
1	G	34	LEU
1	G	52	ALA
1	G	58	PRO
1	B	42	ALA
1	G	33	GLU
1	G	39	ARG
1	G	40	THR
1	G	42	ALA
1	G	43	LYS
1	G	44	ALA
1	E	235	GLU
1	G	22	ALA
1	G	57	SER

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	199/214 (93%)	181 (91%)	18 (9%)	9	6
1	B	188/214 (88%)	178 (95%)	10 (5%)	20	20
1	C	200/214 (94%)	184 (92%)	16 (8%)	11	8
1	D	193/214 (90%)	183 (95%)	10 (5%)	21	20
1	E	169/214 (79%)	153 (90%)	16 (10%)	8	5
1	F	196/214 (92%)	183 (93%)	13 (7%)	15	13
1	G	137/214 (64%)	125 (91%)	12 (9%)	9	6
1	H	181/214 (85%)	168 (93%)	13 (7%)	13	10
All	All	1463/1712 (86%)	1355 (93%)	108 (7%)	13	10

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	27	LEU
1	A	29	GLU
1	A	34	LEU
1	A	35	THR
1	A	47	LEU
1	A	54	VAL
1	A	62	LEU
1	A	99	SER
1	A	144	ARG
1	A	177	LEU
1	A	189	LEU
1	A	234	ASP
1	A	236	ASP
1	A	245	MET
1	A	246	GLU
1	A	270	ARG
1	A	288	LEU
1	B	16	LEU

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Mol	Chain	Res	Type
1	B	83	GLU
1	B	99	SER
1	B	189	LEU
1	B	192	LEU
1	B	194	VAL
1	B	198	GLU
1	B	236	ASP
1	B	245	MET
1	B	279	GLU
1	C	12	THR
1	C	29	GLU
1	C	47	LEU
1	C	54	VAL
1	C	62	LEU
1	C	86	GLU
1	C	90	LYS
1	C	99	SER
1	C	189	LEU
1	C	192	LEU
1	C	211	GLU
1	C	228	GLU
1	C	234	ASP
1	C	236	ASP
1	C	245	MET
1	C	288	LEU
1	D	12	THR
1	D	13	LEU
1	D	47	LEU
1	D	118	ARG
1	D	131	ILE
1	D	192	LEU
1	D	245	MET
1	D	249	VAL
1	D	259	THR
1	D	263	GLU
1	E	12	THR
1	E	21	GLN
1	E	72	LEU
1	E	81	VAL
1	E	99	SER
1	E	118	ARG
1	E	121	THR

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Mol	Chain	Res	Type
1	E	144	ARG
1	E	148	GLU
1	E	153	LEU
1	E	211	GLU
1	E	215	MET
1	E	236	ASP
1	E	245	MET
1	E	246	GLU
1	E	253	ARG
1	F	9	SER
1	F	38	ASN
1	F	73	THR
1	F	99	SER
1	F	135	SER
1	F	158	THR
1	F	171	GLN
1	F	189	LEU
1	F	211	GLU
1	F	236	ASP
1	F	245	MET
1	F	270	ARG
1	F	284	ARG
1	G	19	MET
1	G	21	GLN
1	G	23	MET
1	G	121	THR
1	G	163	ARG
1	G	171	GLN
1	G	175	GLN
1	G	236	ASP
1	G	240	MET
1	G	245	MET
1	G	246	GLU
1	G	288	LEU
1	H	12	THR
1	H	14	ILE
1	H	16	LEU
1	H	21	GLN
1	H	54	VAL
1	H	99	SER
1	H	144	ARG
1	H	158	THR

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Mol	Chain	Res	Type
1	H	171	GLN
1	H	198	GLU
1	H	245	MET
1	H	263	GLU
1	H	284	ARG

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

Mol	Chain	Res	Type
1	A	38	ASN
1	A	75	HIS
1	A	175	GLN
1	A	247	HIS
1	B	75	HIS
1	B	175	GLN
1	B	190	HIS
1	C	175	GLN
1	D	38	ASN
1	D	75	HIS
1	D	175	GLN
1	F	38	ASN
1	F	175	GLN
1	G	75	HIS
1	G	171	GLN
1	G	213	ASN
1	H	38	ASN
1	H	175	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

2 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
2	BOG	F	1000	-	20,20,20	1.25	1 (5%)	25,25,25	1.16	2 (8%)
2	BOG	H	1000	-	20,20,20	1.82	4 (20%)	25,25,25	2.10	5 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BOG	F	1000	-	-	6/11/31/31	0/1/1/1
2	BOG	H	1000	-	-	6/11/31/31	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1000	BOG	O1-C1	6.08	1.50	1.40
2	F	1000	BOG	O1-C1	3.94	1.46	1.40
2	H	1000	BOG	C4-C3	2.91	1.59	1.52
2	H	1000	BOG	O5-C1	2.39	1.48	1.41
2	H	1000	BOG	C3-C2	2.15	1.57	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1000	BOG	O5-C1-O1	7.60	128.00	110.04
2	F	1000	BOG	C6'-C5'-C4'	-2.88	99.80	114.37
2	H	1000	BOG	C6'-C5'-C4'	-2.80	100.23	114.37
2	H	1000	BOG	C3'-C2'-C1'	-2.74	101.55	113.47
2	F	1000	BOG	C3'-C2'-C1'	-2.74	101.56	113.47
2	H	1000	BOG	C1'-O1-C1	2.26	117.54	113.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1000	BOG	O5-C5-C6	2.01	111.43	106.44

There are no chirality outliers.

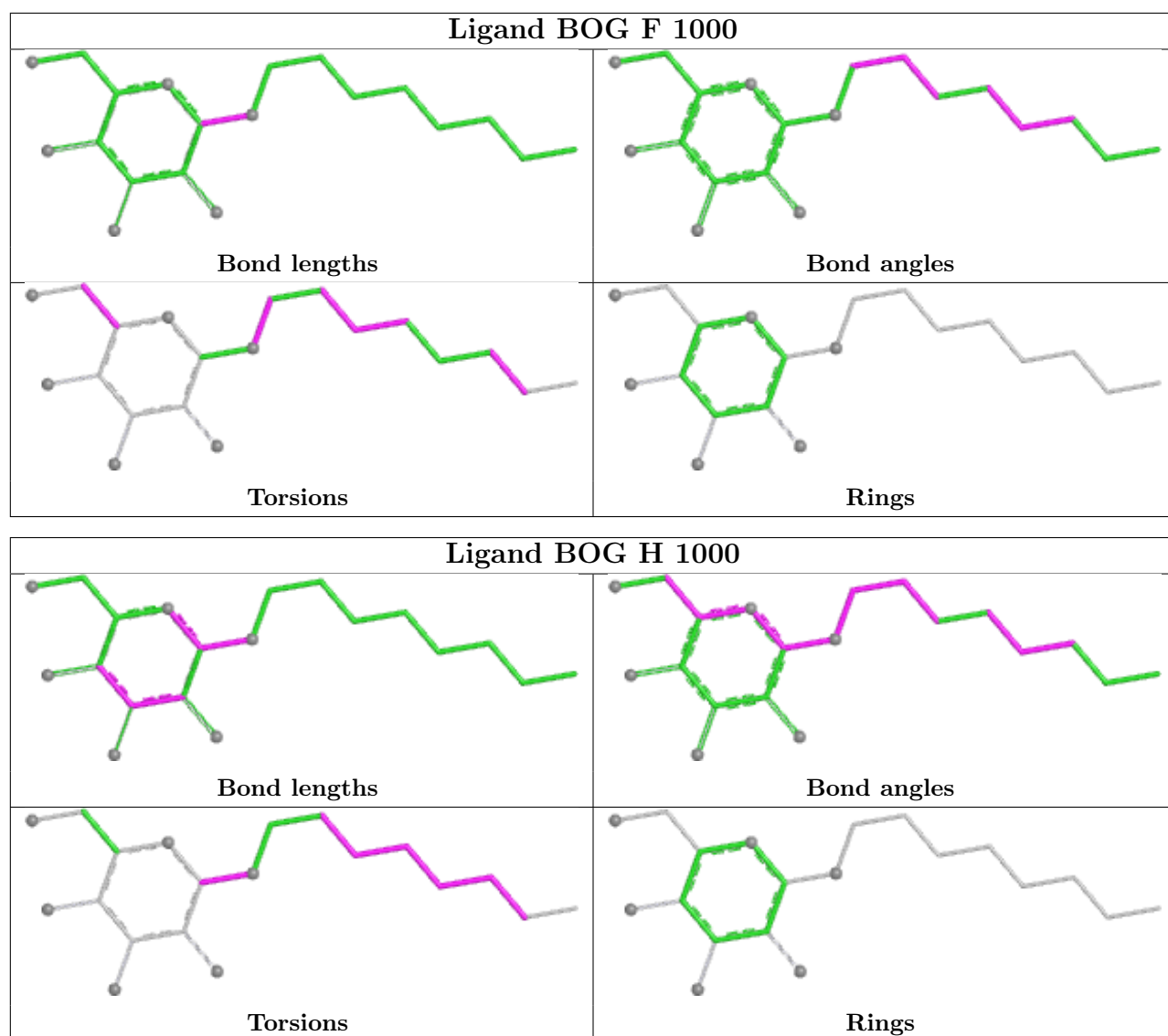
All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	1000	BOG	O5-C1-O1-C1'
2	F	1000	BOG	C2'-C3'-C4'-C5'
2	H	1000	BOG	C4'-C5'-C6'-C7'
2	F	1000	BOG	C1'-C2'-C3'-C4'
2	H	1000	BOG	C1'-C2'-C3'-C4'
2	H	1000	BOG	C2'-C3'-C4'-C5'
2	H	1000	BOG	C5'-C6'-C7'-C8'
2	F	1000	BOG	C5'-C6'-C7'-C8'
2	F	1000	BOG	C4-C5-C6-O6
2	F	1000	BOG	O5-C5-C6-O6
2	F	1000	BOG	C2'-C1'-O1-C1
2	H	1000	BOG	C3'-C4'-C5'-C6'

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.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	F	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	F	13:LEU	C	14:ILE	N	1.01

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	285/293 (97%)	0.54	11 (3%)	43 43	26, 43, 70, 102	0
1	B	282/293 (96%)	0.45	18 (6%)	25 24	26, 45, 76, 102	0
1	C	283/293 (96%)	0.40	7 (2%)	58 59	25, 40, 66, 88	0
1	D	285/293 (97%)	0.21	9 (3%)	50 50	26, 40, 72, 93	0
1	E	267/293 (91%)	0.93	39 (14%)	6 5	25, 56, 96, 117	0
1	F	283/293 (96%)	0.35	12 (4%)	40 39	23, 40, 76, 102	0
1	G	269/293 (91%)	1.43	94 (34%)	1 0	27, 70, 124, 141	0
1	H	278/293 (94%)	1.30	63 (22%)	2 1	28, 56, 92, 109	0
All	All	2232/2344 (95%)	0.69	253 (11%)	10 9	23, 46, 94, 141	0

All (253) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	42	ALA	7.2
1	G	36	VAL	5.6
1	H	212	GLY	5.5
1	H	236	ASP	5.3
1	H	214	ASP	5.2
1	G	27	LEU	4.5
1	G	93	THR	4.5
1	G	119	TYR	4.4
1	G	59	ALA	4.3
1	G	35	THR	4.2
1	E	145	ALA	4.2
1	E	58	PRO	4.2
1	B	235	GLU	4.1
1	G	85	ALA	4.1
1	G	153	LEU	4.1
1	G	118	ARG	4.1

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Mol	Chain	Res	Type	RSRZ
1	E	54	VAL	4.1
1	G	97	LEU	4.0
1	H	27	LEU	4.0
1	H	234	ASP	4.0
1	B	39	ARG	4.0
1	G	37	TRP	4.0
1	E	40	THR	4.0
1	G	54	VAL	3.9
1	B	232	PRO	3.8
1	E	234	ASP	3.7
1	G	88	ASP	3.7
1	G	55	ALA	3.7
1	G	128	ALA	3.7
1	G	121	THR	3.6
1	G	64	ALA	3.6
1	G	234	ASP	3.6
1	G	95	LEU	3.6
1	G	26	ALA	3.6
1	H	131	ILE	3.5
1	A	42	ALA	3.5
1	G	94	ILE	3.5
1	H	47	LEU	3.5
1	G	40	THR	3.5
1	G	57	SER	3.5
1	E	214	ASP	3.5
1	H	52	ALA	3.5
1	H	56	ASP	3.5
1	G	150	ASN	3.4
1	G	32	TYR	3.4
1	G	56	ASP	3.4
1	G	44	ALA	3.4
1	H	39	ARG	3.4
1	H	34	LEU	3.3
1	H	154	LEU	3.3
1	G	76	ALA	3.3
1	G	145	ALA	3.3
1	G	69	LEU	3.3
1	H	235	GLU	3.3
1	G	70	LEU	3.3
1	E	230	ARG	3.3
1	C	110	ALA	3.2
1	G	147	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	G	213	ASN	3.2
1	H	200	VAL	3.2
1	F	235	GLU	3.1
1	E	84	GLY	3.1
1	G	61	ALA	3.1
1	H	42	ALA	3.1
1	G	34	LEU	3.1
1	G	135	SER	3.1
1	A	291	ALA	3.1
1	G	19	MET	3.1
1	E	57	SER	3.0
1	G	105	SER	3.0
1	H	232	PRO	3.0
1	D	232	PRO	3.0
1	E	55	ALA	3.0
1	G	24	ALA	3.0
1	F	39	ARG	3.0
1	H	134	SER	3.0
1	H	256	GLY	3.0
1	G	20	GLY	3.0
1	A	76	ALA	2.9
1	C	59	ALA	2.9
1	E	11	VAL	2.9
1	G	120	VAL	2.9
1	H	54	VAL	2.9
1	H	68	VAL	2.9
1	E	233	GLY	2.9
1	G	212	GLY	2.9
1	C	44	ALA	2.9
1	E	28	LEU	2.9
1	E	34	LEU	2.9
1	H	24	ALA	2.9
1	H	197	ALA	2.9
1	B	44	ALA	2.9
1	G	47	LEU	2.9
1	H	45	GLU	2.9
1	C	43	LYS	2.8
1	G	72	LEU	2.8
1	H	282	PHE	2.8
1	E	62	LEU	2.8
1	B	275	ALA	2.8
1	E	24	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	H	30	ALA	2.8
1	H	55	ALA	2.8
1	D	134	SER	2.8
1	G	134	SER	2.8
1	A	44	ALA	2.7
1	F	234	ASP	2.7
1	G	60	GLU	2.7
1	G	137	PHE	2.7
1	G	156	VAL	2.7
1	A	43	LYS	2.7
1	E	37	TRP	2.7
1	G	154	LEU	2.7
1	H	64	ALA	2.6
1	E	56	ASP	2.6
1	H	36	VAL	2.6
1	H	12	THR	2.6
1	H	49	GLU	2.6
1	B	236	ASP	2.6
1	G	83	GLU	2.6
1	F	11	VAL	2.6
1	F	40	THR	2.6
1	B	198	GLU	2.6
1	C	49	GLU	2.6
1	H	44	ALA	2.6
1	G	111	TRP	2.6
1	G	68	VAL	2.6
1	G	77	VAL	2.6
1	H	129	PRO	2.6
1	E	87	SER	2.6
1	H	28	LEU	2.5
1	H	46	ALA	2.5
1	H	48	ALA	2.5
1	D	234	ASP	2.5
1	A	45	GLU	2.5
1	B	45	GLU	2.5
1	C	76	ALA	2.5
1	G	22	ALA	2.5
1	G	46	ALA	2.5
1	H	152	ALA	2.5
1	H	291	ALA	2.5
1	B	91	GLY	2.5
1	G	115	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	42	ALA	2.5
1	G	127	PRO	2.5
1	E	39	ARG	2.5
1	E	211	GLU	2.4
1	F	9	SER	2.4
1	E	64	ALA	2.4
1	H	26	ALA	2.4
1	A	235	GLU	2.4
1	G	114	GLY	2.4
1	B	135	SER	2.4
1	G	33	GLU	2.4
1	E	121	THR	2.4
1	B	41	LYS	2.4
1	D	136	VAL	2.4
1	E	16	LEU	2.4
1	G	89	LEU	2.4
1	H	87	SER	2.4
1	H	10	PRO	2.3
1	H	58	PRO	2.3
1	E	36	VAL	2.3
1	G	81	VAL	2.3
1	G	157	LEU	2.3
1	E	38	ASN	2.3
1	B	291	ALA	2.3
1	E	85	ALA	2.3
1	G	108	ALA	2.3
1	C	240	MET	2.3
1	D	40	THR	2.3
1	H	69	LEU	2.3
1	A	41	LYS	2.3
1	F	10	PRO	2.3
1	G	79	TYR	2.3
1	A	40	THR	2.3
1	H	93	THR	2.3
1	F	282	PHE	2.3
1	F	232	PRO	2.3
1	G	84	GLY	2.3
1	G	98	GLY	2.3
1	H	25	GLY	2.3
1	H	84	GLY	2.3
1	G	217	MET	2.3
1	E	119	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	42	ALA	2.3
1	E	61	ALA	2.3
1	G	39	ARG	2.3
1	G	71	SER	2.3
1	G	131	ILE	2.3
1	B	194	VAL	2.3
1	H	202	VAL	2.3
1	G	91	GLY	2.2
1	G	144	ARG	2.2
1	H	230	ARG	2.2
1	H	9	SER	2.2
1	H	94	ILE	2.2
1	F	13	LEU	2.2
1	G	143	ASP	2.2
1	B	228	GLU	2.2
1	E	235	GLU	2.2
1	G	29	GLU	2.2
1	G	160	THR	2.2
1	H	135	SER	2.2
1	H	252	SER	2.2
1	E	68	VAL	2.2
1	A	217	MET	2.2
1	G	78	MET	2.2
1	H	122	GLY	2.2
1	B	43	LYS	2.2
1	G	48	ALA	2.2
1	G	110	ALA	2.2
1	H	59	ALA	2.2
1	D	39	ARG	2.2
1	G	62	LEU	2.2
1	G	92	ARG	2.2
1	H	226	VAL	2.2
1	E	67	PRO	2.2
1	E	229	GLY	2.2
1	H	57	SER	2.2
1	G	230	ARG	2.1
1	G	23	MET	2.1
1	H	11	VAL	2.1
1	H	132	GLY	2.1
1	F	42	ALA	2.1
1	E	32	TYR	2.1
1	F	217	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	213	ASN	2.1
1	G	96	ASN	2.1
1	G	25	GLY	2.1
1	H	65	GLY	2.1
1	B	9	SER	2.1
1	G	103	ALA	2.1
1	D	49	GLU	2.1
1	G	233	GLY	2.1
1	H	259	THR	2.1
1	G	30	ALA	2.1
1	E	153	LEU	2.1
1	G	169	LEU	2.1
1	H	95	LEU	2.1
1	A	77	VAL	2.1
1	H	233	GLY	2.1
1	B	292	GLU	2.0
1	G	155	GLU	2.0
1	D	236	ASP	2.0
1	E	89	LEU	2.0
1	G	129	PRO	2.0
1	G	50	ARG	2.0
1	H	249	VAL	2.0
1	H	31	GLY	2.0
1	E	60	GLU	2.0
1	G	149	GLU	2.0
1	E	70	LEU	2.0
1	G	232	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

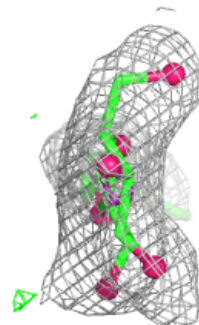
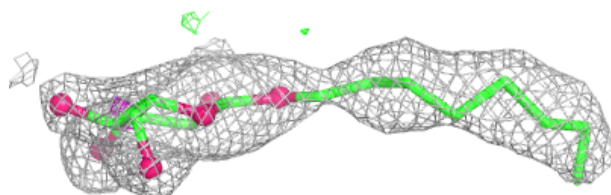
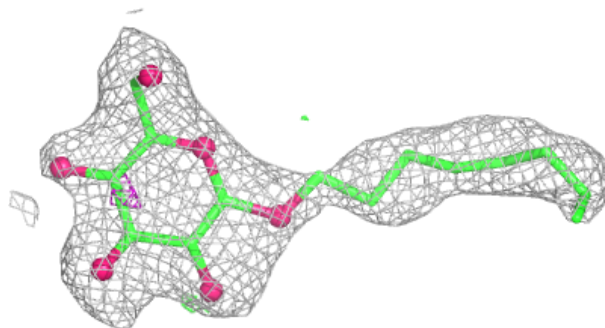
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

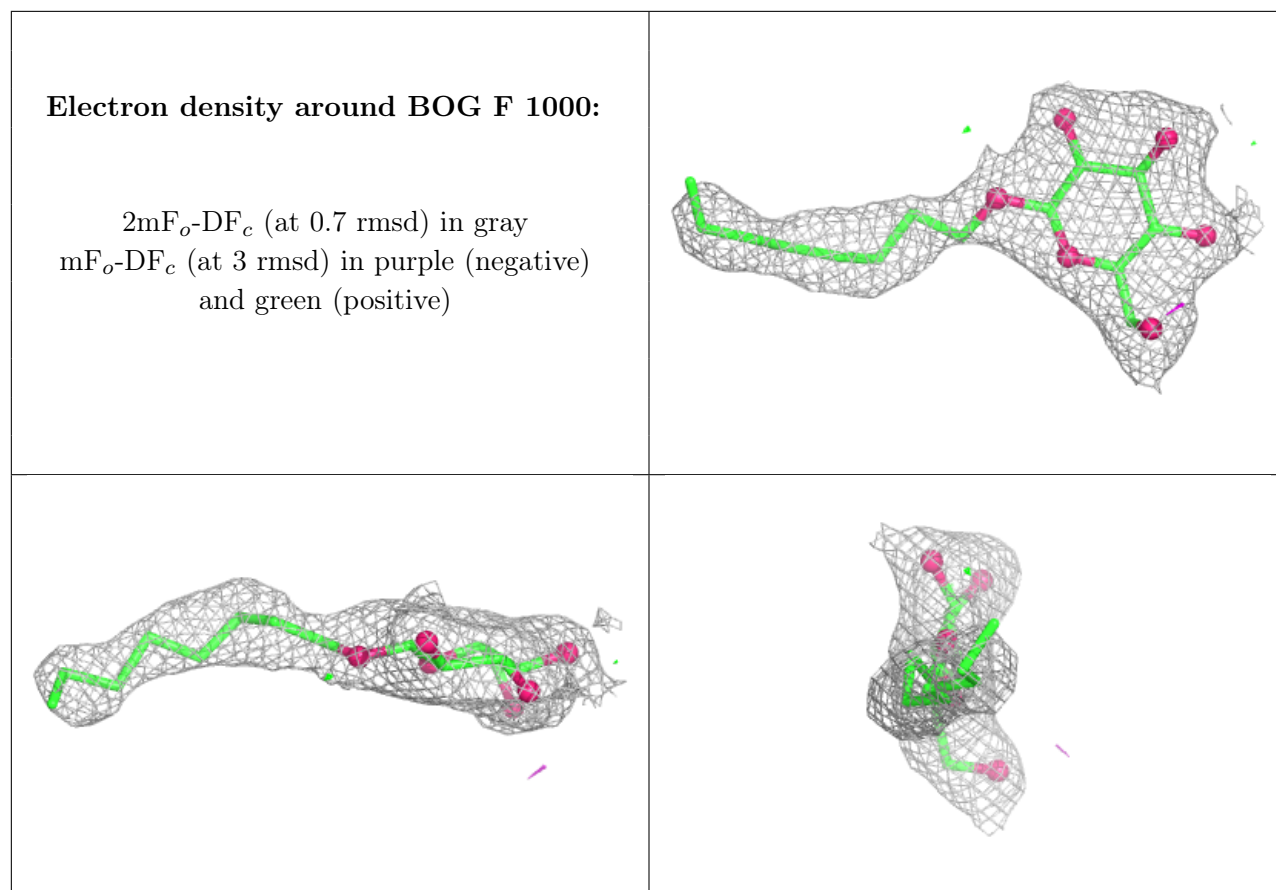
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BOG	H	1000	20/20	0.88	0.15	45,53,68,73	0
2	BOG	F	1000	20/20	0.95	0.10	36,46,69,69	0

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 BOG H 1000:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.