



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 03:50 AM EDT

PDB ID : 1YK3 / pdb_00001yk3
Title : Crystal structure of Rv1347c from Mycobacterium tuberculosis
Authors : Card, G.L.; Peterson, N.A.; Smith, C.A.; Rupp, B.; Schick, B.M.; Baker, E.N.;
TB Structural Genomics Consortium (TBSGC)
Deposited on : 2005-01-16
Resolution : 2.20 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

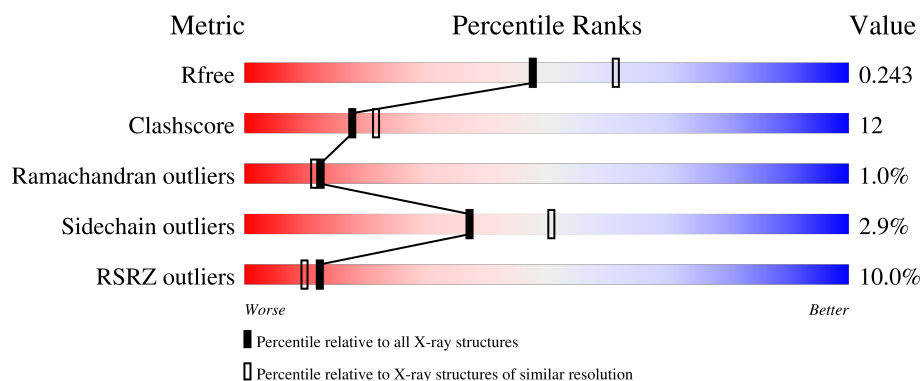
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	210	15% 62% 30% 6%
1	B	210	12% 67% 21% 8%
1	C	210	6% 69% 24% 5%
1	D	210	11% 67% 27% 5%
1	E	210	12% 65% 27% 7%

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Mol	Chain	Length	Quality of chain
1	F	210	4% 78% 18%
1	G	210	3% 77% 18%
1	H	210	12% 65% 28% 6%

2 Entry composition

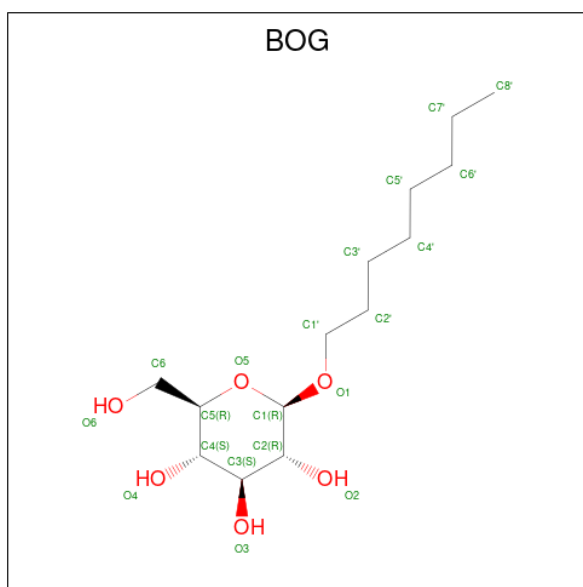
There are 3 unique types of molecules in this entry. The entry contains 13515 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hypothetical protein Rv1347c/MT1389.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	198	Total	C	N	O	S	0	0	0
			1600	1017	292	284	7			
1	B	193	Total	C	N	O	S	0	0	0
			1563	995	285	276	7			
1	C	199	Total	C	N	O	S	0	0	0
			1607	1021	293	286	7			
1	D	200	Total	C	N	O	S	0	0	0
			1612	1024	294	287	7			
1	E	196	Total	C	N	O	S	0	0	0
			1587	1010	290	280	7			
1	F	202	Total	C	N	O	S	0	0	0
			1625	1031	297	290	7			
1	G	201	Total	C	N	O	S	0	0	0
			1618	1027	295	289	7			
1	H	197	Total	C	N	O	S	0	0	0
			1593	1013	291	282	7			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			20	14	6		
2	E	1	Total	C	O	0	0
			20	14	6		
2	G	1	Total	C	O	0	0
			20	14	6		

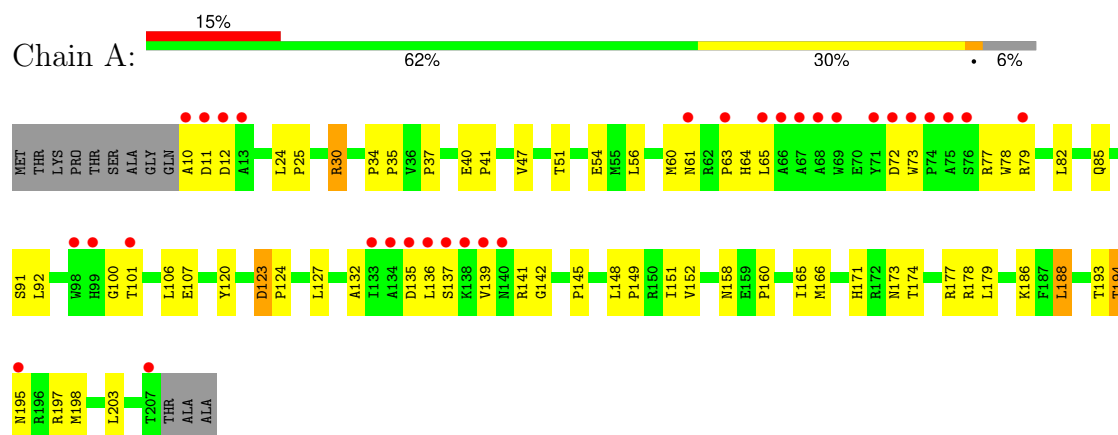
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	51	Total	O	0	0
			51	51		
3	B	54	Total	O	0	0
			54	54		
3	C	83	Total	O	0	0
			83	83		
3	D	84	Total	O	0	0
			84	84		
3	E	59	Total	O	0	0
			59	59		
3	F	136	Total	O	0	0
			136	136		
3	G	121	Total	O	0	0
			121	121		
3	H	62	Total	O	0	0
			62	62		

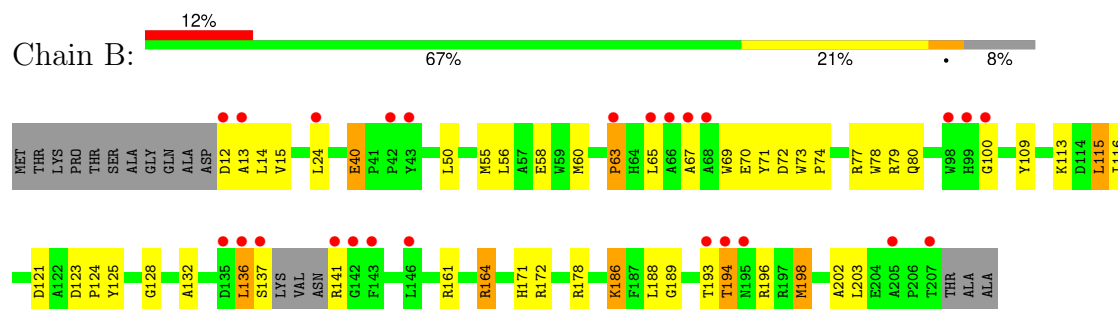
3 Residue-property plots [i](#)

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.

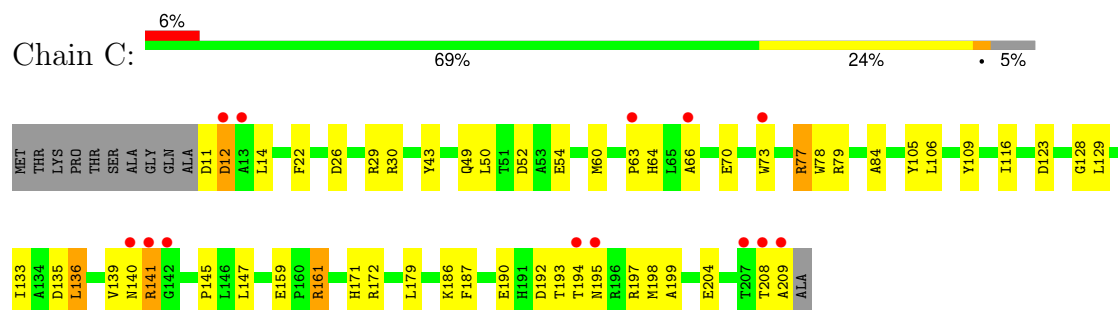
- Molecule 1: Hypothetical protein Rv1347c/MT1389



- Molecule 1: Hypothetical protein Rv1347c/MT1389

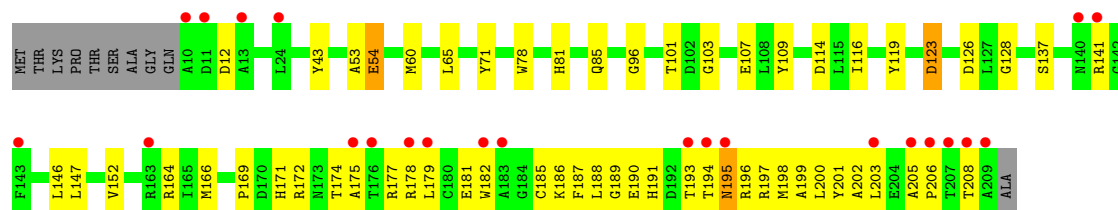


- Molecule 1: Hypothetical protein Rv1347c/MT1389



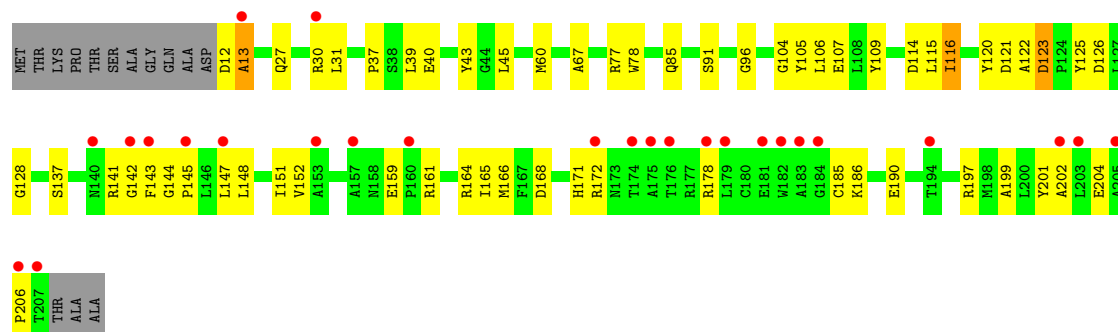
- Molecule 1: Hypothetical protein Rv1347c/MT1389

Chain D: 



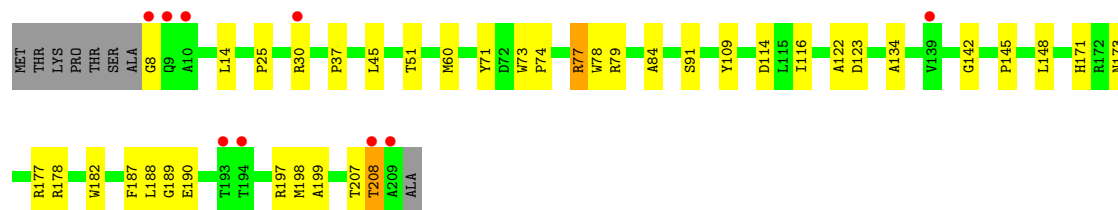
• Molecule 1: Hypothetical protein Rv1347c/MT1389

Chain E: 



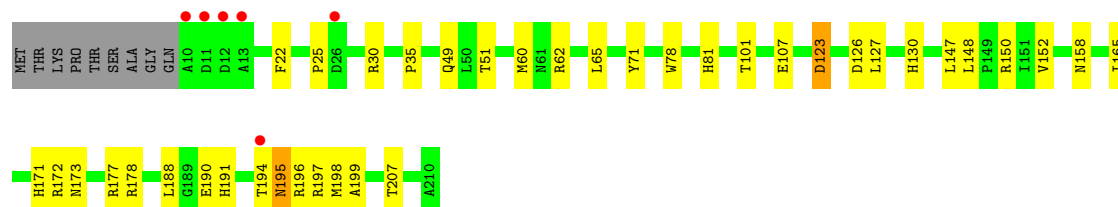
• Molecule 1: Hypothetical protein Rv1347c/MT1389

Chain F: 



• Molecule 1: Hypothetical protein Rv1347c/MT1389

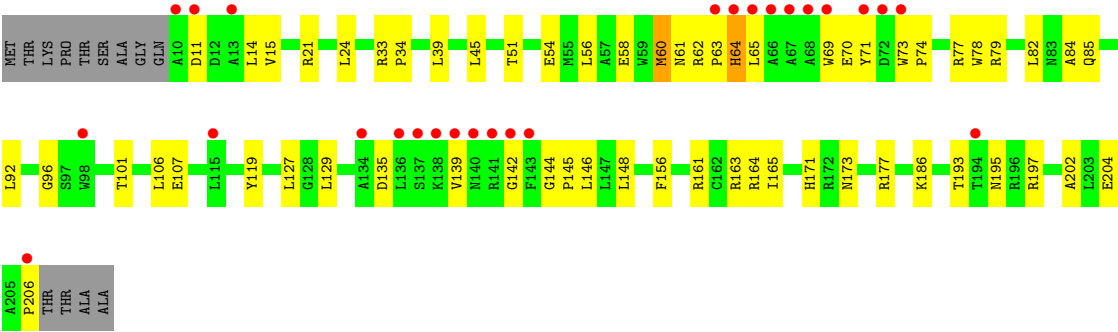
Chain G: 



• Molecule 1: Hypothetical protein Rv1347c/MT1389

Chain H: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.67Å 77.26Å 296.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.96 – 2.20 36.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.9 (36.96-2.20) 96.0 (36.96-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.258 0.218 , 0.243	Depositor DCC
R_{free} test set	1857 reflections (2.08%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13515	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.27% 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

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	0.38	0/1649	0.88	2/2250 (0.1%)
1	B	0.40	0/1611	0.89	1/2197 (0.0%)
1	C	0.41	0/1656	0.90	3/2260 (0.1%)
1	D	0.40	0/1661	0.90	5/2267 (0.2%)
1	E	0.40	0/1636	0.90	2/2232 (0.1%)
1	F	0.44	0/1674	0.92	6/2284 (0.3%)
1	G	0.44	0/1667	0.93	6/2274 (0.3%)
1	H	0.39	0/1642	0.86	1/2240 (0.0%)
All	All	0.41	0/13196	0.90	26/18004 (0.1%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	123	ASP	N-CA-C	-7.46	100.69	110.39
1	C	123	ASP	N-CA-C	-6.46	101.99	110.39
1	F	123	ASP	N-CA-C	-6.44	101.95	110.07
1	D	53	ALA	N-CA-C	6.24	118.60	111.11
1	G	199	ALA	N-CA-C	-6.10	98.46	108.41
1	G	123	ASP	N-CA-C	-6.09	101.88	110.29
1	C	199	ALA	N-CA-C	-6.09	98.48	108.41
1	G	62	ARG	N-CA-C	-6.07	102.12	109.83
1	E	199	ALA	N-CA-C	-5.95	97.42	108.02
1	F	122	ALA	N-CA-C	5.91	118.83	109.96
1	A	123	ASP	N-CA-C	-5.75	101.85	110.24
1	G	130	HIS	N-CA-C	-5.71	100.39	109.59
1	F	71	TYR	N-CA-C	-5.59	103.98	111.87
1	B	189	GLY	N-CA-C	5.50	121.42	112.61
1	H	60	MET	N-CA-C	5.48	118.01	111.71
1	D	123	ASP	N-CA-C	-5.43	102.79	110.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	71	TYR	N-CA-C	-5.43	103.44	111.81
1	A	64	HIS	N-CA-C	5.40	117.16	111.28
1	F	134	ALA	N-CA-C	5.40	116.84	111.07
1	F	25	PRO	N-CA-C	-5.36	103.37	111.41
1	D	12	ASP	N-CA-C	5.32	118.16	107.41
1	F	199	ALA	N-CA-C	-5.32	99.73	108.41
1	G	25	PRO	N-CA-C	-5.30	103.12	111.34
1	C	64	HIS	N-CA-C	5.22	116.97	111.28
1	D	71	TYR	N-CA-C	-5.16	103.86	111.81
1	D	189	GLY	N-CA-C	5.09	119.09	112.68

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	1600	0	1546	48	0
1	B	1563	0	1508	46	0
1	C	1607	0	1553	45	0
1	D	1612	0	1558	45	0
1	E	1587	0	1537	47	0
1	F	1625	0	1569	24	0
1	G	1618	0	1563	28	0
1	H	1593	0	1539	41	0
2	D	20	0	28	3	0
2	E	20	0	28	4	0
2	G	20	0	28	2	0
3	A	51	0	0	2	0
3	B	54	0	0	6	0
3	C	83	0	0	2	0
3	D	84	0	0	1	0
3	E	59	0	0	0	0
3	F	136	0	0	2	0
3	G	121	0	0	4	0
3	H	62	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	13515	0	12457	311	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 12.

All (311) 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:H:60:MET:HE2	1:H:65:LEU:HD13	1.50	0.94
1:H:60:MET:HG3	1:H:78:TRP:CZ2	2.11	0.84
1:B:60:MET:HG3	1:B:78:TRP:CZ2	2.13	0.83
1:B:136:LEU:HD23	1:B:136:LEU:H	1.46	0.80
1:F:60:MET:HG3	1:F:78:TRP:CZ2	2.18	0.79
1:D:190:GLU:HB3	1:D:197:ARG:NH1	2.00	0.77
1:H:106:LEU:HD23	1:H:129:LEU:HD11	1.66	0.77
1:A:158:ASN:HA	1:B:63:PRO:HB3	1.66	0.76
1:B:125:TYR:HE2	1:B:161:ARG:HH21	1.32	0.76
1:C:133:ILE:HB	1:C:139:VAL:HG22	1.67	0.75
1:C:60:MET:HG3	1:C:78:TRP:CZ2	2.22	0.75
1:A:10:ALA:HA	1:E:77:ARG:HH11	1.51	0.74
1:C:116:ILE:HG22	1:C:198:MET:HE1	1.71	0.73
1:D:60:MET:HG3	1:D:78:TRP:CZ2	2.24	0.72
1:G:195:ASN:N	1:G:195:ASN:HD22	1.85	0.72
1:E:85:GLN:HE22	1:E:107:GLU:HG2	1.55	0.71
1:C:43:TYR:CE1	1:C:147:LEU:HD11	2.27	0.70
1:A:60:MET:HG2	1:A:78:TRP:HZ2	1.57	0.69
1:D:172:ARG:NH1	1:D:172:ARG:HB2	2.07	0.69
1:B:116:ILE:HD11	1:B:198:MET:HE2	1.74	0.69
1:G:60:MET:HG3	1:G:78:TRP:CZ2	2.28	0.69
1:D:96:GLY:HA3	2:D:701:BOG:H7'2	1.76	0.68
1:D:190:GLU:HB3	1:D:197:ARG:HH12	1.58	0.68
1:C:136:LEU:HD12	1:C:136:LEU:H	1.59	0.68
1:D:85:GLN:HE22	1:D:107:GLU:HG2	1.59	0.68
1:E:60:MET:HG3	1:E:78:TRP:CZ2	2.29	0.67
1:H:85:GLN:HE22	1:H:107:GLU:HG2	1.58	0.67
1:C:171:HIS:CD2	1:C:197:ARG:HB3	2.29	0.67
1:H:142:GLY:C	1:H:145:PRO:HD2	2.20	0.66
1:E:114:ASP:OD1	1:E:115:LEU:HD23	1.95	0.66
1:D:172:ARG:HB2	1:D:172:ARG:HH11	1.61	0.66
1:H:73:TRP:HB3	1:H:74:PRO:HD2	1.78	0.66
1:C:22:PHE:HB3	1:G:22:PHE:HB3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:173:ASN:O	1:H:177:ARG:HG3	1.96	0.66
1:G:195:ASN:HD22	1:G:195:ASN:H	1.45	0.65
1:G:150:ARG:HB2	3:G:819:HOH:O	1.96	0.65
1:D:171:HIS:CD2	1:D:197:ARG:HB3	2.31	0.65
1:B:171:HIS:CE1	1:B:172:ARG:HG3	2.32	0.65
1:D:116:ILE:HD11	1:D:200:LEU:CD1	2.28	0.64
1:A:100:GLY:HA2	1:D:187:PHE:HB2	1.79	0.64
1:E:123:ASP:HB3	1:E:126:ASP:OD1	1.97	0.64
1:D:175:ALA:HA	1:D:178:ARG:HE	1.62	0.64
1:E:96:GLY:HA3	2:E:702:BOG:H7'2	1.80	0.64
1:H:56:LEU:HD12	1:H:82:LEU:HD11	1.80	0.64
1:H:171:HIS:CD2	1:H:197:ARG:HB3	2.33	0.64
1:C:193:THR:HG21	1:C:198:MET:HE3	1.80	0.63
1:F:8:GLY:HA3	3:F:219:HOH:O	1.97	0.63
1:A:137:SER:O	1:A:141:ARG:HD3	1.98	0.63
1:H:39:LEU:HD12	1:H:45:LEU:HG	1.81	0.63
1:E:164:ARG:HD2	1:E:202:ALA:HB1	1.81	0.63
1:D:190:GLU:CB	1:D:197:ARG:HH12	2.12	0.62
1:G:81:HIS:HD2	3:G:811:HOH:O	1.81	0.62
1:C:73:TRP:HB3	1:C:77:ARG:HB3	1.82	0.62
1:D:164:ARG:HD3	1:D:202:ALA:HB1	1.81	0.62
1:B:50:LEU:HD13	1:B:79:ARG:HE	1.64	0.62
1:B:136:LEU:H	1:B:136:LEU:CD2	2.12	0.62
1:C:77:ARG:HG3	1:C:77:ARG:HH11	1.64	0.62
1:D:174:THR:O	1:D:178:ARG:HG2	1.99	0.62
1:E:105:TYR:O	1:E:106:LEU:HD23	2.01	0.61
1:F:77:ARG:HG2	1:F:77:ARG:HH11	1.64	0.60
1:H:164:ARG:HD2	1:H:202:ALA:HB1	1.83	0.60
1:H:60:MET:CE	1:H:65:LEU:HD13	2.29	0.60
1:H:69:TRP:O	1:H:71:TYR:N	2.31	0.60
1:B:136:LEU:HD23	1:B:136:LEU:N	2.16	0.59
1:E:85:GLN:HG2	1:E:91:SER:HB3	1.84	0.59
1:H:186:LYS:HE2	1:H:204:GLU:CD	2.26	0.59
1:A:37:PRO:HG3	3:A:238:HOH:O	2.03	0.59
1:B:73:TRP:HB3	1:B:74:PRO:HD2	1.85	0.59
1:C:50:LEU:HG	3:C:274:HOH:O	2.00	0.59
1:E:67:ALA:HB2	1:H:161:ARG:HG2	1.84	0.59
1:C:194:THR:O	1:C:195:ASN:HB2	2.02	0.59
1:E:152:VAL:HG13	1:E:165:ILE:HD12	1.83	0.59
1:C:135:ASP:O	1:C:139:VAL:HG23	2.03	0.58
1:H:69:TRP:C	1:H:71:TYR:H	2.11	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:123:ASP:HB3	1:D:126:ASP:OD1	2.02	0.58
1:E:27:GLN:HG3	1:E:30:ARG:HH21	1.68	0.58
1:B:123:ASP:OD1	1:B:124:PRO:HD2	2.03	0.58
1:H:11:ASP:OD1	1:H:79:ARG:NH1	2.36	0.58
1:F:142:GLY:C	1:F:145:PRO:HD2	2.29	0.58
1:F:30:ARG:C	1:F:30:ARG:HD2	2.29	0.58
1:F:77:ARG:HH11	1:F:77:ARG:CG	2.16	0.58
1:H:135:ASP:O	1:H:139:VAL:HG23	2.03	0.57
1:B:13:ALA:HB3	3:B:225:HOH:O	2.03	0.57
1:F:171:HIS:CD2	1:F:197:ARG:HB3	2.39	0.57
1:G:196:ARG:HG2	1:G:196:ARG:HH11	1.67	0.57
1:A:142:GLY:C	1:A:145:PRO:HD2	2.29	0.57
1:C:192:ASP:CG	1:C:197:ARG:HD2	2.30	0.57
1:H:58:GLU:HG2	1:H:62:ARG:NH2	2.19	0.57
1:D:101:THR:HG21	2:D:701:BOG:H2	1.87	0.57
1:C:194:THR:HG22	1:C:195:ASN:ND2	2.19	0.57
1:E:37:PRO:HG2	1:E:45:LEU:HD23	1.86	0.57
1:D:81:HIS:HE1	1:D:107:GLU:OE2	1.87	0.57
1:G:152:VAL:HG13	1:G:165:ILE:HD12	1.87	0.57
1:E:115:LEU:HD23	1:E:116:ILE:N	2.19	0.56
1:B:60:MET:CE	1:B:65:LEU:HD13	2.35	0.56
1:C:159:GLU:OE2	1:C:161:ARG:HD3	2.05	0.56
1:E:116:ILE:HD11	1:E:168:ASP:HB3	1.86	0.56
1:E:166:MET:HE1	1:E:202:ALA:HB2	1.88	0.56
1:D:190:GLU:OE1	1:D:197:ARG:NH1	2.39	0.56
1:G:195:ASN:N	1:G:195:ASN:ND2	2.48	0.56
1:F:91:SER:HB3	1:F:109:TYR:HB3	1.87	0.56
1:A:135:ASP:O	1:A:139:VAL:HG23	2.06	0.55
1:A:56:LEU:HD12	1:A:82:LEU:HD11	1.88	0.55
1:G:147:LEU:HD13	2:G:703:BOG:H7'2	1.89	0.55
1:B:109:TYR:CZ	1:B:128:GLY:HA3	2.41	0.55
1:C:192:ASP:OD1	1:C:197:ARG:HD2	2.06	0.55
1:D:146:LEU:HD23	1:D:146:LEU:O	2.06	0.55
1:E:30:ARG:HG3	1:E:30:ARG:HH11	1.72	0.55
1:E:85:GLN:NE2	1:E:107:GLU:HG2	2.22	0.55
1:A:152:VAL:HG13	1:A:165:ILE:HD12	1.89	0.54
1:D:43:TYR:CE1	1:D:147:LEU:HD11	2.42	0.54
1:B:60:MET:HE2	1:B:65:LEU:HD13	1.89	0.54
1:F:178:ARG:HG3	1:F:182:TRP:CE2	2.42	0.54
1:H:14:LEU:HD11	1:H:84:ALA:HA	1.90	0.53
1:E:137:SER:O	1:E:141:ARG:HG3	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:141:ARG:HG2	1:E:141:ARG:HH11	1.73	0.53
1:H:186:LYS:HE2	1:H:204:GLU:OE2	2.09	0.53
1:D:116:ILE:HD11	1:D:200:LEU:HD11	1.91	0.53
1:A:51:THR:HG21	1:D:119:TYR:CE2	2.44	0.53
1:G:190:GLU:OE1	1:G:197:ARG:NH1	2.42	0.53
1:H:60:MET:HG3	1:H:78:TRP:HZ2	1.71	0.53
1:A:171:HIS:CD2	1:A:197:ARG:HB3	2.44	0.52
1:B:164:ARG:HG3	1:B:202:ALA:HB1	1.90	0.52
1:G:81:HIS:HE1	1:G:107:GLU:OE2	1.92	0.52
1:H:61:ASN:C	1:H:62:ARG:HH11	2.17	0.52
1:H:21:ARG:HG3	3:H:241:HOH:O	2.09	0.52
1:B:193:THR:O	1:B:194:THR:C	2.53	0.52
1:A:30:ARG:HH11	1:A:30:ARG:HB2	1.74	0.52
1:B:77:ARG:HA	1:B:80:GLN:HE21	1.74	0.52
1:H:144:GLY:O	1:H:148:LEU:HD23	2.10	0.52
1:F:37:PRO:HG2	1:F:45:LEU:HD13	1.92	0.52
1:H:127:LEU:O	1:H:165:ILE:HA	2.09	0.51
1:A:41:PRO:HG3	1:D:186:LYS:HE3	1.92	0.51
1:A:186:LYS:HE3	1:A:188:LEU:HD11	1.91	0.51
1:C:136:LEU:H	1:C:136:LEU:CD1	2.23	0.51
1:E:12:ASP:O	1:E:13:ALA:HB2	2.09	0.51
1:A:10:ALA:HA	1:E:77:ARG:NH1	2.23	0.51
1:D:81:HIS:HD2	3:D:741:HOH:O	1.92	0.51
1:G:51:THR:HG23	3:G:758:HOH:O	2.10	0.51
1:A:54:GLU:CD	1:A:54:GLU:H	2.18	0.51
1:B:115:LEU:CD1	1:B:194:THR:HG23	2.41	0.50
1:C:54:GLU:N	1:C:54:GLU:OE2	2.43	0.50
1:D:193:THR:HG22	1:D:196:ARG:O	2.11	0.50
1:D:137:SER:O	1:D:141:ARG:HG2	2.11	0.50
1:A:30:ARG:HB2	1:A:30:ARG:NH1	2.26	0.50
1:B:69:TRP:O	1:B:70:GLU:HB2	2.12	0.50
1:B:60:MET:HG3	1:B:78:TRP:HZ2	1.70	0.49
1:C:105:TYR:C	1:C:106:LEU:HD12	2.37	0.49
1:C:73:TRP:CE3	1:C:77:ARG:HD2	2.47	0.49
1:C:77:ARG:HG3	1:C:77:ARG:NH1	2.27	0.49
1:C:136:LEU:HD12	1:C:136:LEU:N	2.24	0.49
1:G:173:ASN:O	1:G:177:ARG:HG3	2.13	0.49
1:A:160:PRO:HD2	1:B:67:ALA:CB	2.43	0.49
1:A:123:ASP:OD1	1:A:124:PRO:HD2	2.13	0.49
1:G:49:GLN:HA	1:G:49:GLN:NE2	2.27	0.49
1:E:115:LEU:HD23	1:E:116:ILE:H	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:178:ARG:HG3	1:F:182:TRP:CZ2	2.47	0.48
1:B:69:TRP:NE1	1:B:132:ALA:HB2	2.28	0.48
1:E:144:GLY:N	1:E:145:PRO:HD2	2.28	0.48
1:D:172:ARG:HH11	1:D:172:ARG:CB	2.26	0.48
1:B:100:GLY:HA2	1:C:187:PHE:HB2	1.95	0.48
1:C:140:ASN:HB3	1:C:141:ARG:NH2	2.28	0.48
1:D:103:GLY:HA3	2:D:701:BOG:H2'2	1.95	0.48
1:C:60:MET:HG3	1:C:78:TRP:HZ2	1.71	0.48
1:E:43:TYR:CE1	1:E:147:LEU:HD11	2.49	0.48
1:G:127:LEU:O	1:G:165:ILE:HA	2.14	0.48
1:D:175:ALA:O	1:D:179:LEU:HG	2.14	0.47
1:H:146:LEU:HG	3:H:272:HOH:O	2.12	0.47
1:D:114:ASP:HB3	1:D:116:ILE:HG22	1.96	0.47
1:E:39:LEU:HD21	1:E:151:ILE:HG12	1.95	0.47
1:F:77:ARG:CG	1:F:77:ARG:NH1	2.75	0.47
1:H:11:ASP:CG	1:H:79:ARG:HH12	2.21	0.47
1:B:55:MET:O	1:B:58:GLU:HB3	2.13	0.47
1:B:73:TRP:HB3	1:B:77:ARG:HB3	1.95	0.47
1:C:14:LEU:HD11	1:C:84:ALA:HA	1.95	0.47
1:E:144:GLY:O	1:E:148:LEU:HD13	2.15	0.47
1:H:15:VAL:HG22	3:H:253:HOH:O	2.15	0.47
1:B:115:LEU:HD13	3:B:262:HOH:O	2.14	0.47
1:C:29:ARG:HD3	3:C:257:HOH:O	2.14	0.47
1:E:190:GLU:HB3	1:E:197:ARG:NH1	2.30	0.47
1:H:69:TRP:C	1:H:71:TYR:N	2.73	0.47
1:B:65:LEU:HD22	1:B:132:ALA:HB1	1.97	0.47
1:F:60:MET:HG3	1:F:78:TRP:HZ2	1.78	0.46
1:G:171:HIS:NE2	1:G:172:ARG:HG3	2.30	0.46
1:F:188:LEU:HD12	1:F:188:LEU:N	2.29	0.46
1:C:30:ARG:HG3	1:C:30:ARG:HH11	1.79	0.46
1:E:186:LYS:HZ1	1:E:204:GLU:HG3	1.80	0.46
1:A:60:MET:HE3	1:A:78:TRP:CH2	2.50	0.46
1:F:8:GLY:N	1:F:79:ARG:HD3	2.31	0.46
1:B:69:TRP:HE1	1:B:132:ALA:HB2	1.80	0.46
1:D:116:ILE:HD11	1:D:200:LEU:HD13	1.95	0.46
1:E:178:ARG:HA	1:E:178:ARG:HD3	1.75	0.46
1:B:203:LEU:C	1:B:203:LEU:HD23	2.41	0.46
1:F:114:ASP:OD1	1:F:116:ILE:HG23	2.16	0.46
1:A:194:THR:HG23	1:A:195:ASN:N	2.31	0.46
1:E:120:TYR:CE1	1:E:122:ALA:HA	2.50	0.46
1:A:40:GLU:HA	1:A:40:GLU:OE1	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:TRP:O	1:A:82:LEU:HG	2.15	0.45
1:E:109:TYR:CZ	1:E:128:GLY:HA3	2.51	0.45
1:A:174:THR:HG22	1:A:178:ARG:NE	2.31	0.45
1:D:188:LEU:O	1:D:191:HIS:HE1	1.98	0.45
1:H:119:TYR:CZ	1:H:193:THR:HG22	2.51	0.45
1:C:50:LEU:HB3	1:C:79:ARG:NH1	2.30	0.45
1:E:30:ARG:HG3	1:E:30:ARG:NH1	2.30	0.45
1:C:73:TRP:HB3	1:C:77:ARG:CB	2.46	0.45
1:A:51:THR:OG1	1:D:191:HIS:HD2	2.00	0.45
1:A:73:TRP:CE3	1:A:77:ARG:HG3	2.52	0.45
1:B:186:LYS:HE2	3:B:243:HOH:O	2.16	0.45
1:D:152:VAL:HG11	1:D:203:LEU:HD23	1.99	0.45
1:F:187:PHE:C	1:F:188:LEU:HD12	2.41	0.45
1:H:62:ARG:HB3	1:H:63:PRO:HD2	1.99	0.45
1:H:62:ARG:HA	1:H:62:ARG:HD3	1.81	0.45
1:B:63:PRO:HB2	3:B:256:HOH:O	2.16	0.45
1:H:34:PRO:O	1:H:92:LEU:HD22	2.17	0.45
1:B:12:ASP:CG	1:B:13:ALA:H	2.25	0.45
1:D:178:ARG:HD2	1:D:182:TRP:CH2	2.51	0.44
1:E:106:LEU:HD11	1:E:151:ILE:HD13	1.98	0.44
1:C:194:THR:O	1:C:195:ASN:CB	2.65	0.44
1:A:47:VAL:HG23	3:A:233:HOH:O	2.17	0.44
1:D:205:ALA:HB1	1:D:206:PRO:HD2	2.00	0.44
1:A:127:LEU:O	1:A:165:ILE:HA	2.17	0.44
1:A:136:LEU:HD23	1:A:136:LEU:O	2.16	0.44
1:D:185:CYS:HB3	1:D:201:TYR:HB3	2.00	0.44
1:F:207:THR:O	1:F:208:THR:C	2.60	0.44
1:E:171:HIS:HB3	1:E:197:ARG:HG2	1.98	0.44
1:A:61:ASN:OD1	1:A:72:ASP:HB2	2.17	0.44
1:B:56:LEU:HD13	1:B:78:TRP:CZ3	2.53	0.44
1:A:65:LEU:HD22	1:A:132:ALA:HB1	1.98	0.44
1:C:109:TYR:CZ	1:C:128:GLY:HA3	2.53	0.44
1:G:49:GLN:HA	1:G:49:GLN:HE21	1.82	0.43
1:G:123:ASP:HB3	1:G:126:ASP:OD1	2.17	0.43
1:B:124:PRO:HB2	1:B:125:TYR:CE1	2.53	0.43
1:F:145:PRO:HD3	3:F:280:HOH:O	2.18	0.43
1:C:186:LYS:HE2	1:C:204:GLU:HG3	2.00	0.43
1:E:143:PHE:CZ	2:E:702:BOG:H4'2	2.54	0.43
1:H:156:PHE:CD1	1:H:206:PRO:HD3	2.54	0.43
1:A:106:LEU:HD21	1:A:151:ILE:HD13	2.01	0.43
1:C:208:THR:O	1:C:209:ALA:HB3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:187:PHE:CZ	1:F:189:GLY:HA2	2.54	0.43
1:A:24:LEU:HA	1:A:25:PRO:HD3	1.88	0.43
1:B:71:TYR:O	1:B:73:TRP:N	2.48	0.43
1:C:49:GLN:O	1:C:52:ASP:HB2	2.18	0.43
1:D:54:GLU:CD	1:D:54:GLU:H	2.26	0.43
1:D:190:GLU:HA	1:D:198:MET:O	2.18	0.43
1:G:196:ARG:HH11	1:G:196:ARG:CG	2.32	0.43
1:A:120:TYR:CZ	1:A:166:MET:HE2	2.54	0.43
1:D:166:MET:HE1	1:D:188:LEU:HD12	2.00	0.43
1:A:173:ASN:O	1:A:177:ARG:HG3	2.18	0.43
1:B:40:GLU:OE1	1:B:40:GLU:HA	2.19	0.43
1:E:143:PHE:CE1	2:E:702:BOG:H4'2	2.54	0.43
1:F:190:GLU:HA	1:F:198:MET:O	2.19	0.43
1:E:27:GLN:HG3	1:E:30:ARG:NH2	2.33	0.43
1:A:85:GLN:NE2	1:A:91:SER:OG	2.47	0.42
1:F:173:ASN:O	1:F:177:ARG:HG3	2.19	0.42
1:G:49:GLN:NE2	3:G:783:HOH:O	2.51	0.42
1:B:125:TYR:N	1:B:125:TYR:CD1	2.87	0.42
1:A:85:GLN:HE22	1:A:107:GLU:HG2	1.83	0.42
1:B:116:ILE:CD1	1:B:198:MET:HE2	2.46	0.42
1:B:137:SER:O	1:B:141:ARG:HB2	2.19	0.42
1:E:185:CYS:HB3	1:E:201:TYR:HB3	2.01	0.42
1:G:101:THR:HG21	2:G:703:BOG:H2	2.00	0.42
1:A:85:GLN:HG2	1:A:91:SER:HB3	2.01	0.42
1:A:137:SER:HA	1:A:141:ARG:NH1	2.35	0.42
1:B:178:ARG:HG2	3:B:263:HOH:O	2.19	0.42
1:E:31:LEU:HD21	1:E:125:TYR:CE2	2.54	0.42
1:G:30:ARG:HG2	1:G:30:ARG:HH11	1.85	0.42
1:A:65:LEU:CD2	1:A:132:ALA:HB1	2.50	0.42
1:A:148:LEU:HB2	1:A:149:PRO:HD3	2.02	0.42
1:G:35:PRO:HB2	1:G:158:ASN:HB3	2.01	0.42
1:G:178:ARG:HG3	1:G:178:ARG:HH11	1.85	0.42
1:C:145:PRO:HB3	1:C:179:LEU:HD21	2.02	0.42
1:H:45:LEU:HD23	1:H:96:GLY:HA2	2.02	0.42
1:B:100:GLY:HA2	1:C:187:PHE:CB	2.50	0.42
1:E:104:GLY:HA3	2:E:702:BOG:H7'1	2.02	0.42
1:A:60:MET:HG2	1:A:78:TRP:CZ2	2.46	0.41
1:C:11:ASP:O	1:C:12:ASP:C	2.63	0.41
1:F:14:LEU:HD11	1:F:84:ALA:HA	2.02	0.41
1:B:15:VAL:HG13	3:B:212:HOH:O	2.19	0.41
1:B:186:LYS:HD2	1:B:188:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:73:TRP:HB3	1:F:74:PRO:HD2	2.01	0.41
1:B:193:THR:HG22	1:B:196:ARG:O	2.20	0.41
1:C:106:LEU:HD23	1:C:129:LEU:HD11	2.03	0.41
1:D:169:PRO:HD2	1:D:199:ALA:O	2.21	0.41
1:H:142:GLY:O	1:H:145:PRO:HD2	2.21	0.41
1:H:64:HIS:CD2	1:H:64:HIS:N	2.88	0.41
1:A:79:ARG:HH11	1:A:79:ARG:HG2	1.85	0.41
1:B:24:LEU:HD13	1:B:113:LYS:HG2	2.01	0.41
1:C:73:TRP:CB	1:C:77:ARG:HB3	2.49	0.41
1:D:177:ARG:O	1:D:181:GLU:HG3	2.21	0.41
1:E:172:ARG:HD2	1:H:163:ARG:NE	2.36	0.41
1:A:34:PRO:HA	1:A:35:PRO:HD3	1.84	0.41
1:A:60:MET:HE3	1:A:78:TRP:CZ2	2.56	0.41
1:E:186:LYS:HG3	1:E:204:GLU:OE2	2.21	0.41
1:G:188:LEU:O	1:G:191:HIS:HE1	2.03	0.41
1:H:73:TRP:HB3	1:H:77:ARG:HB3	2.03	0.41
1:C:73:TRP:CD2	1:C:77:ARG:HD2	2.56	0.41
1:D:65:LEU:HD23	1:D:65:LEU:HA	1.94	0.41
1:D:195:ASN:C	1:D:195:ASN:HD22	2.25	0.41
1:E:40:GLU:HA	1:E:40:GLU:OE2	2.20	0.41
1:E:159:GLU:OE1	1:E:161:ARG:NH2	2.53	0.41
1:A:34:PRO:O	1:A:92:LEU:HD22	2.21	0.41
1:G:171:HIS:CD2	1:G:172:ARG:HG3	2.56	0.41
1:H:73:TRP:HB3	1:H:74:PRO:CD	2.49	0.41
1:A:193:THR:O	1:A:194:THR:C	2.65	0.40
1:C:66:ALA:O	1:C:70:GLU:HA	2.21	0.40
1:E:142:GLY:O	1:E:145:PRO:HG2	2.22	0.40
1:E:190:GLU:HB3	1:E:197:ARG:HH11	1.86	0.40
1:G:65:LEU:HA	1:G:65:LEU:HD23	1.86	0.40
1:D:109:TYR:CZ	1:D:128:GLY:HA3	2.57	0.40
1:C:171:HIS:CD2	1:C:172:ARG:HG3	2.56	0.40
1:C:190:GLU:HA	1:C:198:MET:O	2.22	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	196/210 (93%)	187 (95%)	6 (3%)	3 (2%)	8	6
1	B	189/210 (90%)	183 (97%)	4 (2%)	2 (1%)	11	10
1	C	197/210 (94%)	188 (95%)	7 (4%)	2 (1%)	12	11
1	D	198/210 (94%)	192 (97%)	3 (2%)	3 (2%)	8	6
1	E	194/210 (92%)	184 (95%)	8 (4%)	2 (1%)	12	11
1	F	200/210 (95%)	190 (95%)	9 (4%)	1 (0%)	24	27
1	G	199/210 (95%)	194 (98%)	4 (2%)	1 (0%)	24	27
1	H	195/210 (93%)	190 (97%)	3 (2%)	2 (1%)	12	11
All	All	1568/1680 (93%)	1508 (96%)	44 (3%)	16 (1%)	12	11

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	194	THR
1	E	13	ALA
1	F	208	THR
1	A	11	ASP
1	D	195	ASN
1	H	70	GLU
1	H	195	ASN
1	A	194	THR
1	G	194	THR
1	B	72	ASP
1	C	12	ASP
1	D	208	THR
1	D	194	THR
1	E	206	PRO
1	A	63	PRO
1	C	63	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	162/170 (95%)	155 (96%)	7 (4%)	26	35
1	B	158/170 (93%)	149 (94%)	9 (6%)	18	23
1	C	163/170 (96%)	158 (97%)	5 (3%)	35	48
1	D	163/170 (96%)	162 (99%)	1 (1%)	78	89
1	E	161/170 (95%)	159 (99%)	2 (1%)	63	78
1	F	164/170 (96%)	161 (98%)	3 (2%)	51	68
1	G	163/170 (96%)	159 (98%)	4 (2%)	42	56
1	H	161/170 (95%)	155 (96%)	6 (4%)	30	41
All	All	1295/1360 (95%)	1258 (97%)	37 (3%)	37	51

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ASP
1	A	30	ARG
1	A	101	THR
1	A	179	LEU
1	A	188	LEU
1	A	198	MET
1	A	203	LEU
1	B	14	LEU
1	B	40	GLU
1	B	63	PRO
1	B	115	LEU
1	B	121	ASP
1	B	136	LEU
1	B	164	ARG
1	B	186	LYS
1	B	198	MET
1	C	26	ASP
1	C	77	ARG
1	C	136	LEU
1	C	141	ARG
1	C	161	ARG
1	D	54	GLU
1	E	116	ILE
1	E	121	ASP

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Mol	Chain	Res	Type
1	F	51	THR
1	F	77	ARG
1	F	148	LEU
1	G	148	LEU
1	G	195	ASN
1	G	198	MET
1	G	207	THR
1	H	24	LEU
1	H	33	ARG
1	H	51	THR
1	H	54	GLU
1	H	64	HIS
1	H	101	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	80	GLN
1	A	81	HIS
1	A	83	ASN
1	A	85	GLN
1	A	130	HIS
1	A	140	ASN
1	A	171	HIS
1	B	80	GLN
1	B	83	ASN
1	B	85	GLN
1	B	99	HIS
1	B	118	HIS
1	B	158	ASN
1	B	171	HIS
1	C	27	GLN
1	C	85	GLN
1	C	118	HIS
1	C	140	ASN
1	C	158	ASN
1	C	195	ASN
1	D	64	HIS
1	D	81	HIS
1	D	83	ASN
1	D	85	GLN

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Mol	Chain	Res	Type
1	D	99	HIS
1	D	118	HIS
1	D	171	HIS
1	D	191	HIS
1	E	83	ASN
1	E	85	GLN
1	E	99	HIS
1	E	158	ASN
1	F	27	GLN
1	F	83	ASN
1	F	85	GLN
1	F	118	HIS
1	F	140	ASN
1	F	158	ASN
1	F	171	HIS
1	F	191	HIS
1	G	49	GLN
1	G	61	ASN
1	G	81	HIS
1	G	85	GLN
1	G	191	HIS
1	G	195	ASN
1	H	49	GLN
1	H	64	HIS
1	H	83	ASN
1	H	85	GLN
1	H	171	HIS

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

3 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	E	702	-	20,20,20	1.21	3 (15%)	25,25,25	0.63	0
2	BOG	G	703	-	20,20,20	1.08	2 (10%)	25,25,25	0.64	0
2	BOG	D	701	-	20,20,20	1.18	3 (15%)	25,25,25	0.62	0

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	E	702	-	-	3/11/31/31	0/1/1/1
2	BOG	G	703	-	-	0/11/31/31	0/1/1/1
2	BOG	D	701	-	-	3/11/31/31	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	702	BOG	O5-C1	3.02	1.49	1.41
2	D	701	BOG	O5-C1	2.98	1.49	1.41
2	G	703	BOG	O5-C1	2.89	1.49	1.41
2	E	702	BOG	C4-C5	2.45	1.58	1.53
2	D	701	BOG	C4-C5	2.32	1.58	1.53
2	D	701	BOG	O1-C1	2.13	1.43	1.40
2	G	703	BOG	C4-C5	2.06	1.57	1.53
2	E	702	BOG	O1-C1	2.03	1.43	1.40

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

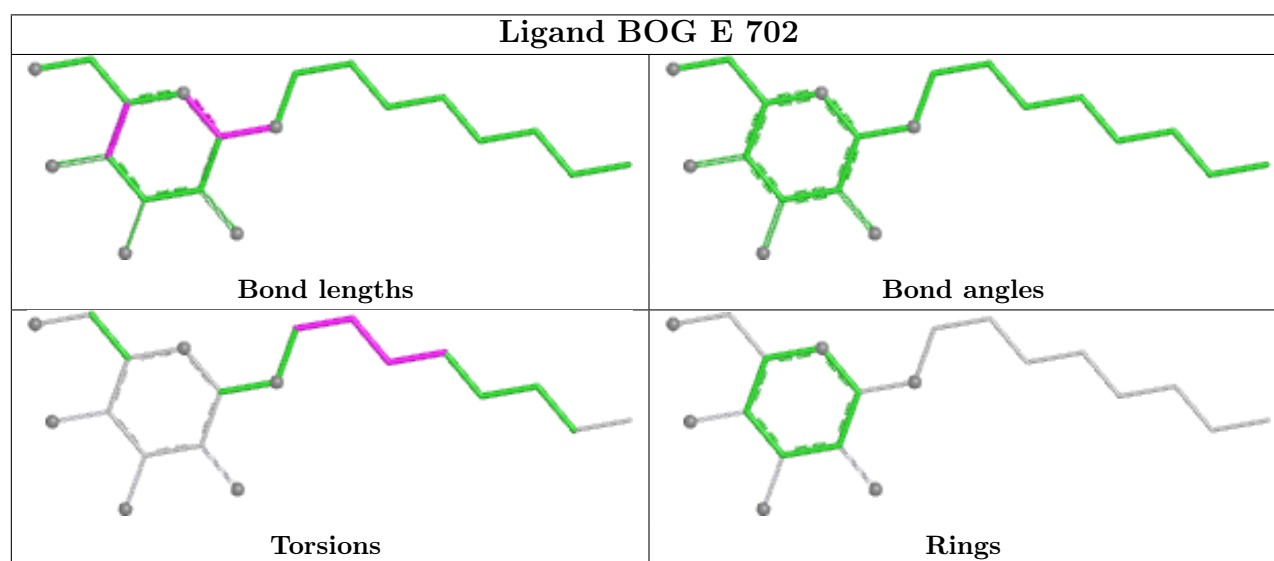
Mol	Chain	Res	Type	Atoms
2	D	701	BOG	O1-C1'-C2'-C3'
2	E	702	BOG	O1-C1'-C2'-C3'
2	D	701	BOG	C2'-C3'-C4'-C5'
2	E	702	BOG	C1'-C2'-C3'-C4'
2	E	702	BOG	C2'-C3'-C4'-C5'
2	D	701	BOG	C1'-C2'-C3'-C4'

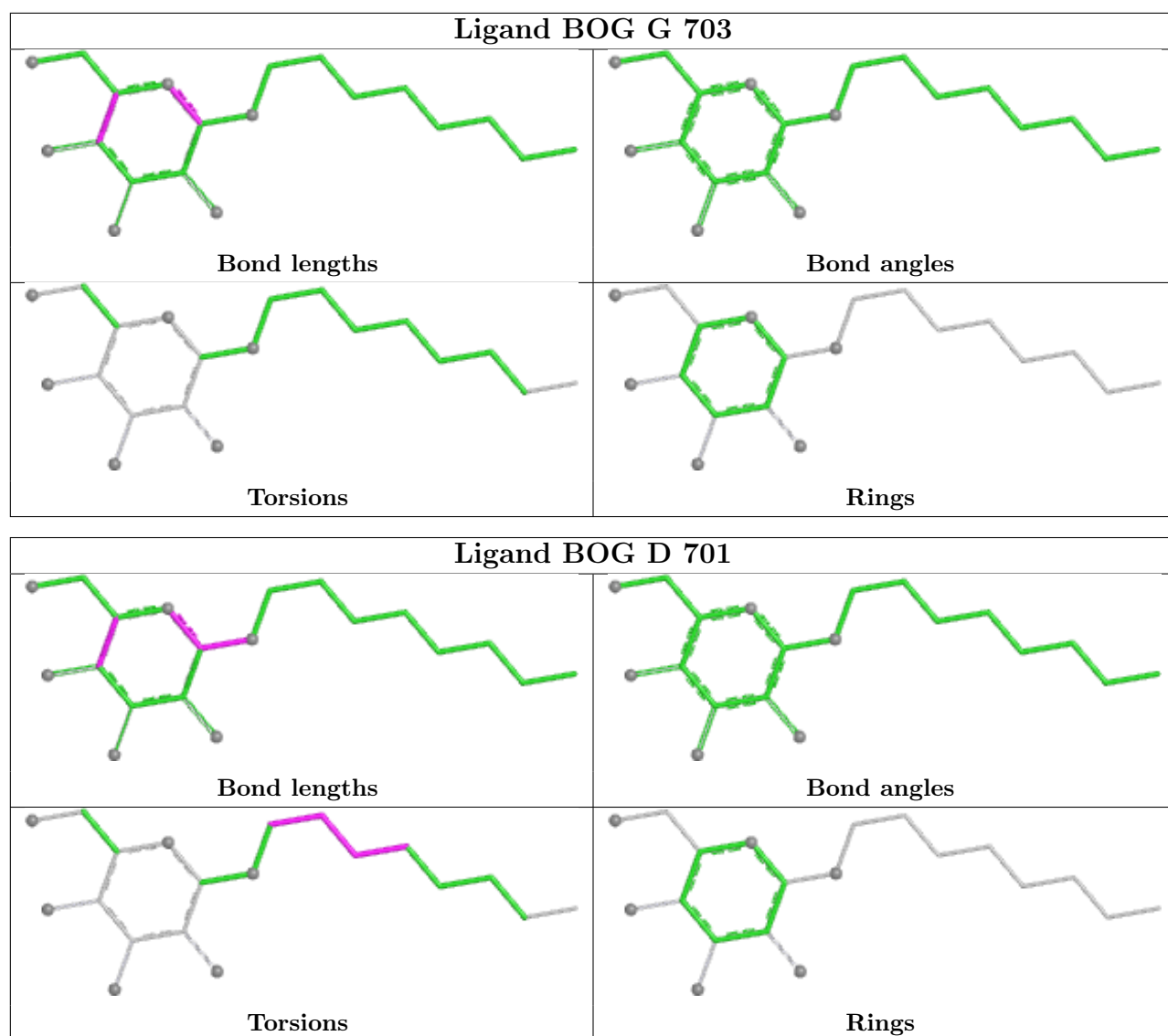
There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	702	BOG	4	0
2	G	703	BOG	2	0
2	D	701	BOG	3	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	198/210 (94%)	0.70	31 (15%) 5 4	23, 37, 76, 82	0
1	B	193/210 (91%)	0.74	25 (12%) 7 5	21, 38, 61, 77	0
1	C	199/210 (94%)	0.32	13 (6%) 25 22	16, 31, 65, 74	0
1	D	200/210 (95%)	0.47	23 (11%) 9 7	18, 32, 59, 79	0
1	E	196/210 (93%)	0.68	26 (13%) 7 5	19, 36, 61, 76	0
1	F	202/210 (96%)	-0.12	9 (4%) 38 35	12, 24, 48, 84	0
1	G	201/210 (95%)	-0.14	6 (2%) 52 49	13, 25, 43, 71	0
1	H	197/210 (93%)	0.84	26 (13%) 7 5	24, 39, 81, 90	0
All	All	1586/1680 (94%)	0.43	159 (10%) 12 10	12, 32, 65, 90	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	8	GLY	5.1
1	H	66	ALA	4.9
1	A	68	ALA	4.6
1	B	143	PHE	4.6
1	F	209	ALA	4.5
1	D	10	ALA	4.5
1	D	208	THR	4.4
1	H	13	ALA	4.4
1	G	12	ASP	4.4
1	E	140	ASN	4.3
1	H	10	ALA	4.3
1	F	10	ALA	4.2
1	A	12	ASP	4.2
1	A	66	ALA	4.1
1	B	135	ASP	4.1
1	C	209	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	139	VAL	4.0
1	B	13	ALA	3.9
1	H	68	ALA	3.9
1	A	71	TYR	3.8
1	A	10	ALA	3.7
1	B	207	THR	3.6
1	H	139	VAL	3.6
1	A	136	LEU	3.5
1	A	13	ALA	3.5
1	D	11	ASP	3.5
1	H	69	TRP	3.5
1	B	12	ASP	3.5
1	E	205	ALA	3.4
1	H	140	ASN	3.4
1	E	142	GLY	3.4
1	E	207	THR	3.4
1	F	208	THR	3.4
1	B	67	ALA	3.4
1	D	209	ALA	3.4
1	H	67	ALA	3.4
1	H	142	GLY	3.4
1	B	66	ALA	3.3
1	H	65	LEU	3.3
1	H	63	PRO	3.3
1	B	142	GLY	3.3
1	A	69	TRP	3.3
1	A	137	SER	3.3
1	A	11	ASP	3.2
1	H	71	TYR	3.2
1	A	135	ASP	3.2
1	A	195	ASN	3.2
1	H	143	PHE	3.1
1	B	194	THR	3.1
1	D	179	LEU	3.1
1	A	75	ALA	3.1
1	A	73	TRP	3.0
1	H	98	TRP	3.0
1	H	137	SER	3.0
1	C	142	GLY	3.0
1	F	194	THR	3.0
1	E	13	ALA	3.0
1	G	194	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	182	TRP	3.0
1	G	10	ALA	2.9
1	E	183	ALA	2.9
1	D	193	THR	2.9
1	E	174	THR	2.9
1	G	11	ASP	2.9
1	A	99	HIS	2.8
1	E	179	LEU	2.8
1	E	184	GLY	2.8
1	B	193	THR	2.7
1	D	13	ALA	2.7
1	F	139	VAL	2.7
1	A	133	ILE	2.7
1	D	178	ARG	2.7
1	B	63	PRO	2.7
1	E	143	PHE	2.7
1	C	140	ASN	2.6
1	H	11	ASP	2.6
1	B	99	HIS	2.6
1	H	64	HIS	2.6
1	C	13	ALA	2.6
1	E	182	TRP	2.6
1	A	140	ASN	2.6
1	A	74	PRO	2.6
1	B	195	ASN	2.6
1	D	207	THR	2.5
1	D	203	LEU	2.5
1	D	140	ASN	2.5
1	G	13	ALA	2.5
1	C	63	PRO	2.5
1	E	172	ARG	2.5
1	H	136	LEU	2.5
1	D	205	ALA	2.5
1	D	195	ASN	2.5
1	E	153	ALA	2.5
1	A	101	THR	2.4
1	C	208	THR	2.4
1	B	98	TRP	2.4
1	H	73	TRP	2.4
1	A	67	ALA	2.4
1	C	66	ALA	2.4
1	C	73	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	195	ASN	2.4
1	E	206	PRO	2.4
1	H	206	PRO	2.4
1	H	194	THR	2.4
1	A	61	ASN	2.4
1	B	42	PRO	2.4
1	D	206	PRO	2.4
1	E	175	ALA	2.4
1	E	176	THR	2.4
1	B	65	LEU	2.4
1	B	141	ARG	2.4
1	C	194	THR	2.3
1	C	207	THR	2.3
1	E	147	LEU	2.3
1	E	202	ALA	2.3
1	B	136	LEU	2.3
1	C	12	ASP	2.3
1	D	183	ALA	2.3
1	H	115	LEU	2.3
1	A	72	ASP	2.3
1	E	157	ALA	2.3
1	H	134	ALA	2.3
1	A	138	LYS	2.3
1	H	141	ARG	2.2
1	D	175	ALA	2.2
1	A	65	LEU	2.2
1	D	141	ARG	2.2
1	E	181	GLU	2.2
1	A	134	ALA	2.2
1	A	207	THR	2.2
1	D	176	THR	2.2
1	C	141	ARG	2.2
1	F	9	GLN	2.2
1	B	146	LEU	2.2
1	A	76	SER	2.2
1	E	194	THR	2.2
1	B	24	LEU	2.1
1	B	205	ALA	2.1
1	E	178	ARG	2.1
1	D	24	LEU	2.1
1	H	138	LYS	2.1
1	A	98	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	100	GLY	2.1
1	E	30	ARG	2.1
1	B	137	SER	2.1
1	B	68	ALA	2.1
1	E	203	LEU	2.1
1	H	72	ASP	2.1
1	E	160	PRO	2.1
1	D	194	THR	2.0
1	F	193	THR	2.0
1	B	43	TYR	2.0
1	A	79	ARG	2.0
1	D	143	PHE	2.0
1	G	26	ASP	2.0
1	A	63	PRO	2.0
1	D	163	ARG	2.0
1	E	145	PRO	2.0
1	F	30	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

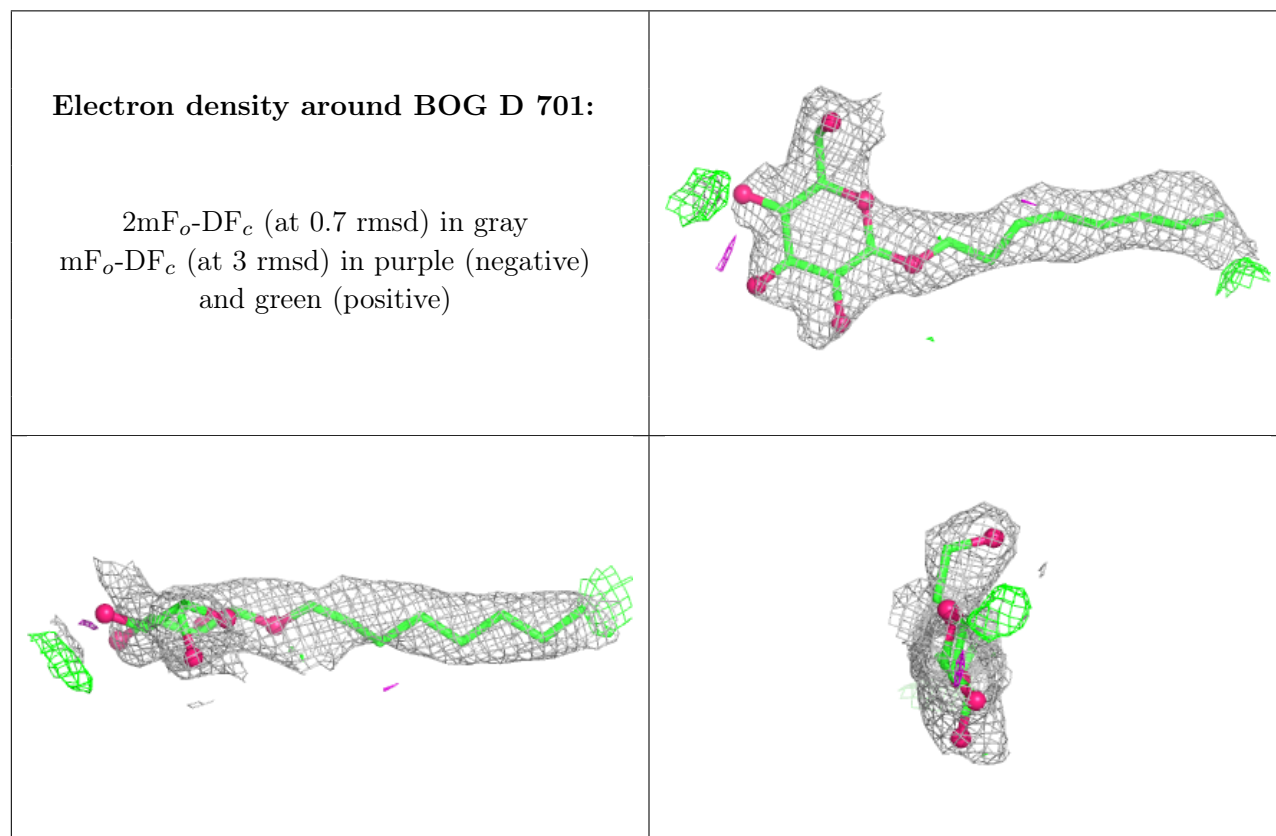
6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BOG	D	701	20/20	0.77	0.18	51,60,63,66	0
2	BOG	E	702	20/20	0.80	0.15	52,61,63,63	0
2	BOG	G	703	20/20	0.81	0.15	50,58,61,62	0

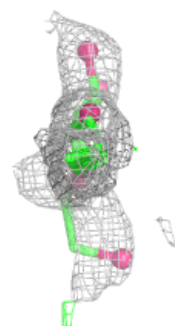
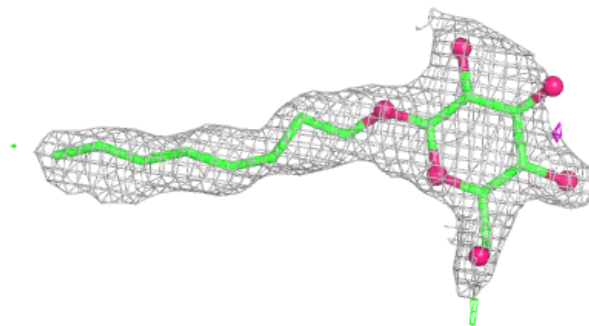
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

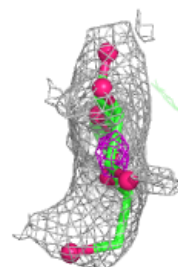
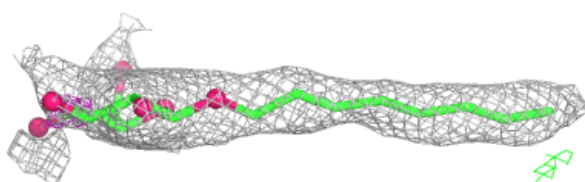
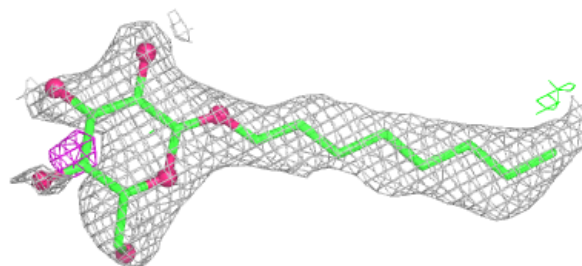


Electron density around BOG E 702:

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

**Electron density around BOG G 703:**

$2mF_o-DF_c$ (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.