



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 11:25 pm BST

PDB ID : 3AUJ  
Title : Structure of diol dehydratase complexed with glycerol  
Authors : Yamanishi, M.; Kinoshita, K.; Fukuoka, M.; Shibata, T.; Tobimatsu, T.; Toraya, T.  
Deposited on : 2011-02-07  
Resolution : 2.10 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

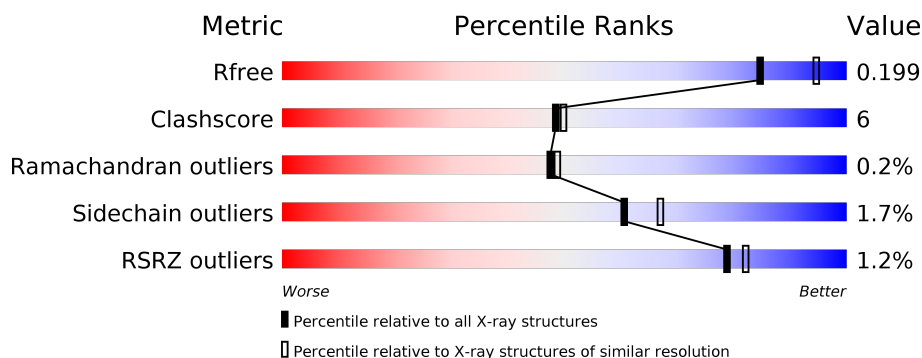
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	554	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>89%</span> <span>10% •</span> </div> </div>
1	L	554	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 97%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>84%</span> <span>14% ••</span> </div> </div>
2	B	224	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 12%, green 86%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>68%</span> <span>12%</span> <span>20%</span> </div> </div>
2	E	224	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 9%, green 89%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>71%</span> <span>9%</span> <span>20%</span> </div> </div>
3	G	173	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 85%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>68%</span> <span>13%</span> <span>20%</span> </div> </div>
3	M	173	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 11%, green 88%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>68%</span> <span>11% •</span> <span>20%</span> </div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14964 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 Diol dehydrase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	0	0
			4210	2625	728	828	29			
1	L	551	Total	C	N	O	S	0	0	0
			4201	2620	727	825	29			

- Molecule 2 is a protein called Diol dehydrase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	179	Total	C	N	O	S	0	0	0
			1367	865	245	255	2			
2	E	180	Total	C	N	O	S	0	0	0
			1375	869	246	258	2			

- Molecule 3 is a protein called Diol dehydrase gamma subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	139	Total	C	N	O	S	0	0	0
			1110	690	200	217	3			
3	M	138	Total	C	N	O	S	0	0	0
			1104	687	199	215	3			

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

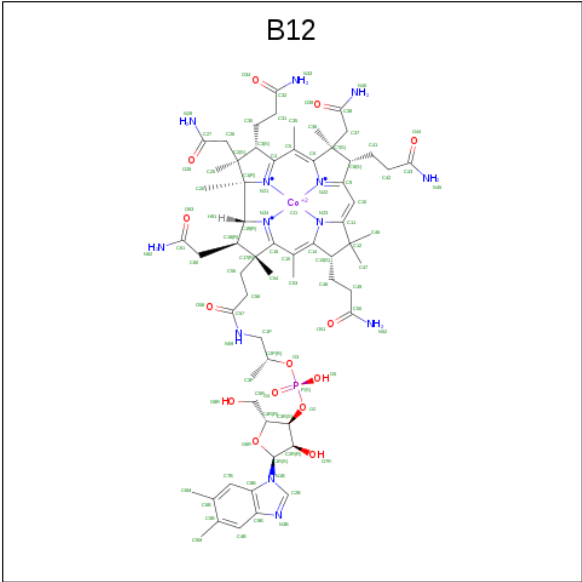
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		
5	L	1	Total	Ca	0	0
			1	1		

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			5	4	1		
6	B	1	Total	O	P	0	0
			5	4	1		
6	G	1	Total	O	P	0	0
			5	4	1		
6	L	1	Total	O	P	0	0
			5	4	1		
6	E	1	Total	O	P	0	0
			5	4	1		
6	M	1	Total	O	P	0	0
			5	4	1		

- Molecule 7 is COBALAMIN (three-letter code: B12) (formula:  $C_{62}H_{89}CoN_{13}O_{14}P$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
7	B	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		
7	E	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		

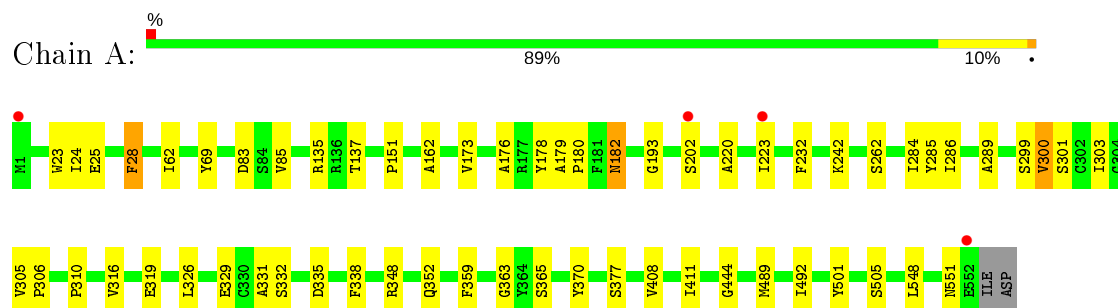
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	417	Total	O	0	0
			417	417		
8	B	158	Total	O	0	0
			158	158		
8	G	116	Total	O	0	0
			116	116		
8	L	393	Total	O	0	0
			393	393		
8	E	178	Total	O	0	0
			178	178		
8	M	109	Total	O	0	0
			109	109		

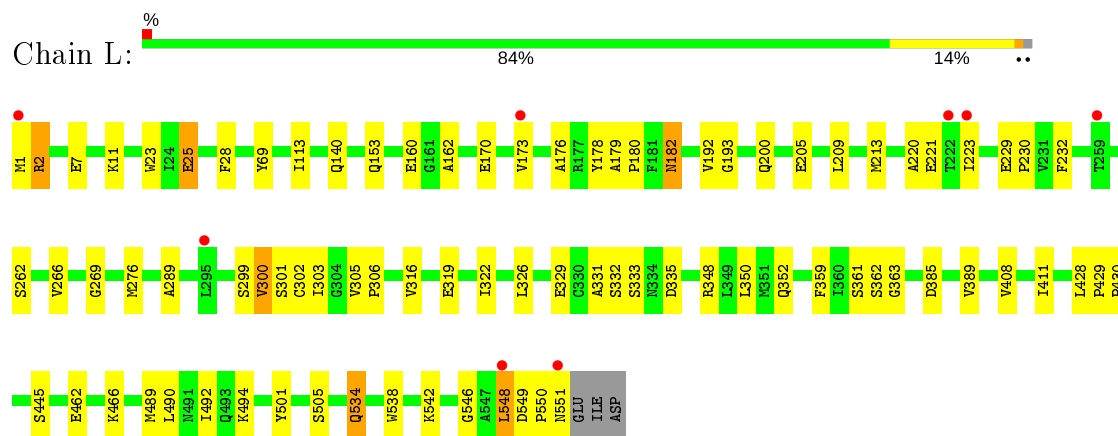
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

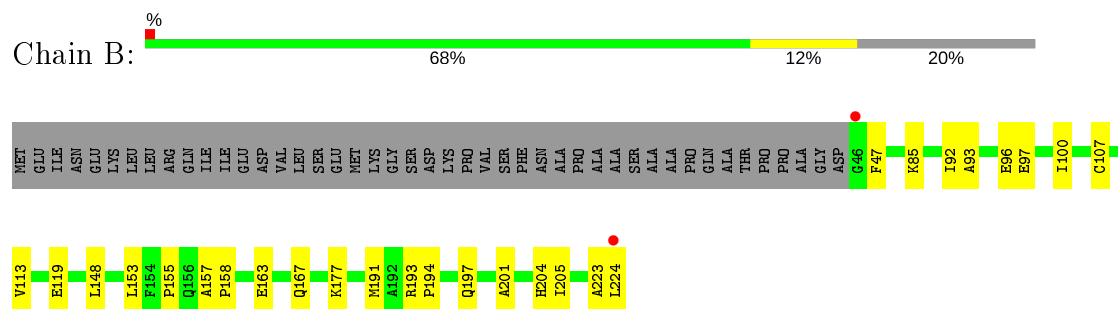
- Molecule 1: Diol dehydrase alpha subunit



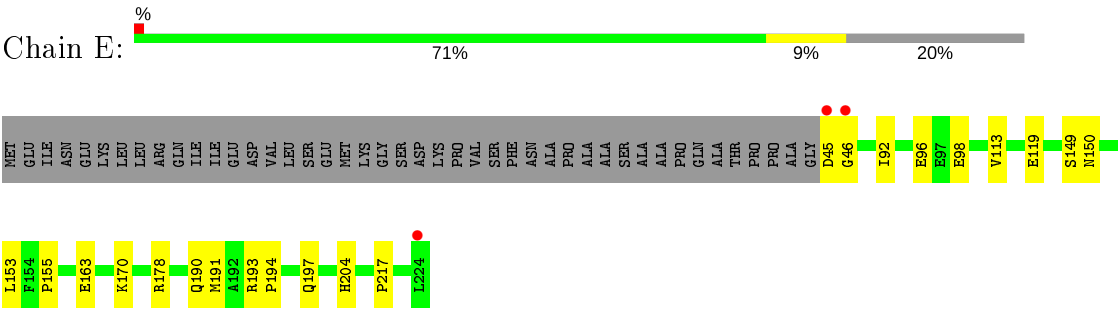
- Molecule 1: Diol dehydrase alpha subunit



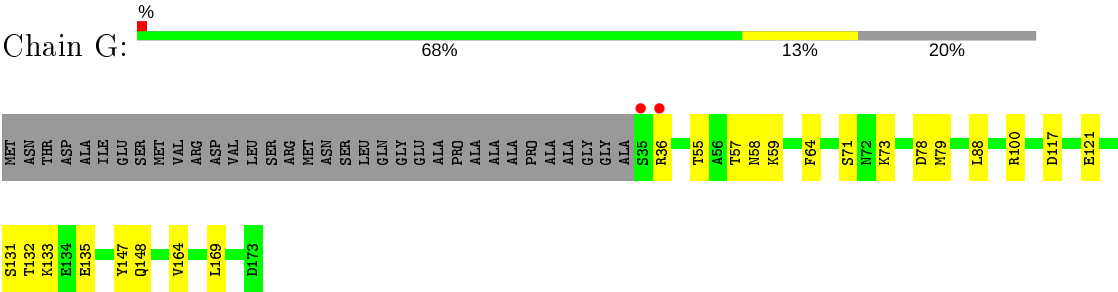
- Molecule 2: Diol dehydrase beta subunit



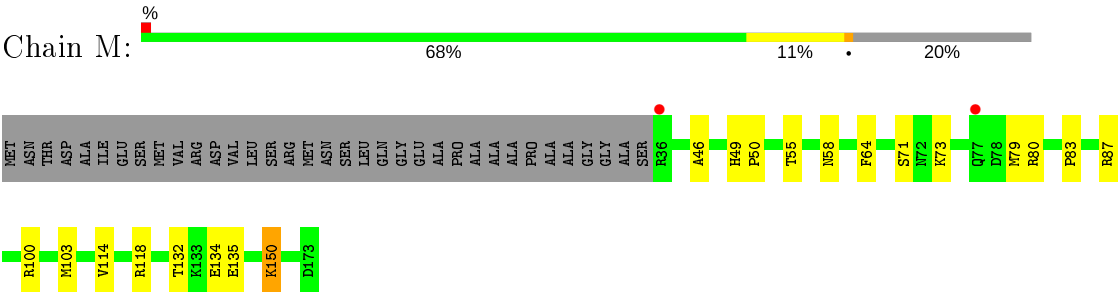
- Molecule 2: Diol dehydrase beta subunit



• Molecule 3: Diol dehydrase gamma subunit



• Molecule 3: Diol dehydrase gamma subunit





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	190.61 Å   196.80 Å   77.36 Å 90.00°   95.03°   90.00°	Depositor
Resolution (Å)	40.34 – 2.10 40.34 – 2.10	Depositor EDS
% Data completeness (in resolution range)	82.8 (40.34-2.10) 81.8 (40.34-2.10)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.81 (at 2.10 Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.186   ,   0.191 0.175   ,   0.199	Depositor DCC
$R_{free}$ test set	15640 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.7	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14964	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CA, B12, PO4

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	0.31	0/4282	0.60	0/5799
1	L	0.31	0/4273	0.60	1/5787 (0.0%)
2	B	0.30	0/1389	0.58	0/1879
2	E	0.30	0/1397	0.58	0/1890
3	G	0.29	0/1125	0.52	0/1519
3	M	0.29	0/1119	0.52	0/1511
All	All	0.31	0/13585	0.58	1/18385 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	L	269	GLY	N-CA-C	5.22	126.14	113.10

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4210	0	4146	38	0
1	L	4201	0	4140	58	0
2	B	1367	0	1419	18	0
2	E	1375	0	1423	15	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1110	0	1119	13	0
3	M	1104	0	1114	16	0
4	A	6	0	6	0	0
4	L	6	0	6	0	0
5	A	1	0	0	0	0
5	L	1	0	0	0	0
6	A	5	0	0	0	0
6	B	5	0	0	0	0
6	E	5	0	0	0	0
6	G	5	0	0	0	0
6	L	5	0	0	0	0
6	M	5	0	0	0	0
7	B	91	0	88	6	0
7	E	91	0	88	5	0
8	A	417	0	0	4	0
8	B	158	0	0	3	0
8	E	178	0	0	1	0
8	G	116	0	0	0	0
8	L	393	0	0	4	0
8	M	109	0	0	1	0
All	All	14964	0	13549	162	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 6.

All (162) 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:L:2:ARG:HB3	1:L:2:ARG:HH11	1.19	1.04
1:A:24:ILE:H	1:A:24:ILE:HD12	1.36	0.90
3:G:71:SER:OG	3:G:73:LYS:HG3	1.83	0.79
3:G:36:ARG:HD2	3:G:88:LEU:HD21	1.66	0.77
1:L:489:MET:O	1:L:492:ILE:HG22	1.87	0.74
1:A:489:MET:O	1:A:492:ILE:HG22	1.88	0.73
3:M:71:SER:OG	3:M:73:LYS:HG2	1.92	0.69
1:L:69:TYR:O	3:M:100:ARG:HD2	1.92	0.69
7:B:1601:B12:H552	7:B:1601:B12:H531	1.74	0.69
3:M:114:VAL:HG13	3:M:118:ARG:HG2	1.76	0.67
3:M:46:ALA:O	3:M:50:PRO:HG3	1.95	0.67
3:M:100:ARG:HH22	3:M:103:MET:HG2	1.58	0.67
7:E:2601:B12:H531	7:E:2601:B12:H552	1.76	0.66

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:107:CYS:HB3	2:B:119:GLU:OE1	1.96	0.66
1:L:140:GLN:OE1	1:L:361:SER:HB2	1.95	0.66
7:B:1601:B12:H351	7:B:1601:B12:H362	1.78	0.66
1:A:24:ILE:H	1:A:24:ILE:CD1	2.09	0.66
1:A:24:ILE:HD12	1:A:24:ILE:N	2.11	0.65
3:G:117:ASP:O	3:G:121:GLU:HG3	1.96	0.65
1:L:2:ARG:CB	1:L:2:ARG:HH11	2.04	0.62
1:L:179:ALA:HB3	1:L:180:PRO:HD3	1.81	0.61
7:E:2601:B12:H362	7:E:2601:B12:H351	1.81	0.61
1:L:408:VAL:O	1:L:411:ILE:HG22	2.01	0.61
1:A:25:GLU:H	1:A:25:GLU:CD	2.04	0.60
3:M:132:THR:OG1	3:M:135:GLU:HG3	2.01	0.60
1:L:2:ARG:HB3	1:L:2:ARG:NH1	2.03	0.60
3:M:100:ARG:NH2	3:M:103:MET:HG2	2.18	0.59
1:L:548:LEU:HD13	1:L:548:LEU:H	1.66	0.59
3:M:83:PRO:O	3:M:87:ARG:HG3	2.02	0.59
1:L:331:ALA:HA	1:L:359:PHE:HB2	1.86	0.58
3:G:132:THR:OG1	3:G:135:GLU:HG3	2.04	0.58
1:L:162:ALA:HB3	1:L:193:GLY:HA3	1.86	0.57
2:E:191:MET:C	2:E:194:PRO:HD2	2.25	0.57
1:A:173:VAL:HG21	1:A:176:ALA:HA	1.85	0.57
2:B:191:MET:C	2:B:194:PRO:HD2	2.25	0.57
1:L:262:SER:HA	8:L:629:HOH:O	2.04	0.57
1:L:192:VAL:O	1:L:411:ILE:HD11	2.05	0.57
3:G:164:VAL:HG22	3:G:169:LEU:HD11	1.86	0.56
1:L:23:TRP:CD1	1:L:25:GLU:HG2	2.40	0.56
1:A:329:GLU:OE2	1:A:505:SER:HA	2.05	0.56
2:B:193:ARG:HB3	2:B:194:PRO:HD3	1.87	0.56
1:L:546:GLY:O	1:L:548:LEU:HD13	2.06	0.56
1:L:113:ILE:HD13	1:L:326:LEU:HD21	1.89	0.55
1:A:332:SER:HB2	1:A:352:GLN:HG2	1.88	0.55
1:A:262:SER:HA	8:A:591:HOH:O	2.07	0.54
3:G:64:PHE:CE1	3:G:79:MET:HG2	2.43	0.53
2:B:93:ALA:O	2:B:97:GLU:HG3	2.08	0.53
1:A:69:TYR:HB2	1:A:289:ALA:HB1	1.90	0.53
1:L:490:LEU:O	1:L:494:LYS:HG3	2.09	0.53
1:L:301:SER:HB3	7:E:2601:B12:H532	1.89	0.53
2:B:223:ALA:CB	2:B:224:LEU:HD12	2.39	0.53
1:L:173:VAL:HG21	1:L:176:ALA:HA	1.91	0.52
1:L:329:GLU:OE2	1:L:505:SER:HA	2.08	0.52
1:L:549:ASP:C	1:L:551:ASN:H	2.12	0.52

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:ASN:N	1:A:182:ASN:HD22	2.08	0.52
1:L:23:TRP:HB3	1:L:28:PHE:HB2	1.92	0.52
1:A:301:SER:HB3	7:B:1601:B12:H532	1.92	0.52
3:G:55:THR:HB	3:G:78:ASP:O	2.10	0.52
2:E:98:GLU:OE1	2:E:170:LYS:HE2	2.10	0.52
3:M:114:VAL:CG1	3:M:118:ARG:HG2	2.40	0.52
3:M:64:PHE:CE1	3:M:79:MET:HG2	2.45	0.52
1:L:305:VAL:HB	1:L:306:PRO:CD	2.40	0.51
2:B:148:LEU:HD13	7:B:1601:B12:H411	1.92	0.51
1:L:332:SER:HB2	1:L:352:GLN:HG2	1.93	0.51
2:E:193:ARG:HB3	2:E:194:PRO:HD3	1.94	0.50
1:L:299:SER:OG	1:L:303:ILE:HA	2.12	0.50
1:L:550:PRO:O	1:L:551:ASN:CG	2.50	0.50
1:A:69:TYR:O	3:G:100:ARG:HD2	2.11	0.50
1:L:322:ILE:O	1:L:326:LEU:HD23	2.11	0.49
3:M:83:PRO:HB2	3:M:87:ARG:NH1	2.27	0.49
1:A:299:SER:OG	1:A:303:ILE:HA	2.12	0.49
2:E:119:GLU:CG	2:E:217:PRO:HG3	2.42	0.49
1:L:69:TYR:HB2	1:L:289:ALA:HB1	1.95	0.49
1:L:205:GLU:HG2	8:L:814:HOH:O	2.13	0.49
1:A:370:TYR:OH	1:A:444:GLY:HA3	2.13	0.48
1:L:220:ALA:HB1	1:L:223:ILE:HD11	1.95	0.48
3:M:80:ARG:HG2	8:M:1253:HOH:O	2.13	0.48
1:L:7:GLU:O	1:L:11:LYS:HG3	2.13	0.48
1:A:179:ALA:HB3	1:A:180:PRO:HD3	1.95	0.48
2:E:92:ILE:O	2:E:96:GLU:HG3	2.13	0.48
2:E:153:LEU:HG	2:E:155:PRO:HD3	1.96	0.48
1:A:284:ILE:HG22	1:A:326:LEU:HD12	1.96	0.47
2:B:223:ALA:HB1	2:B:224:LEU:HD12	1.96	0.47
1:A:300:VAL:HG12	1:A:301:SER:N	2.29	0.47
3:G:55:THR:HG21	3:G:64:PHE:CZ	2.50	0.47
2:B:193:ARG:HG3	2:B:197:GLN:HB2	1.97	0.47
1:L:182:ASN:N	1:L:182:ASN:HD22	2.13	0.47
1:A:285:TYR:CE2	1:A:326:LEU:HD11	2.50	0.47
1:A:83:ASP:OD2	1:A:85:VAL:HB	2.15	0.47
2:E:119:GLU:HG2	2:E:217:PRO:HG3	1.96	0.47
1:A:316:VAL:O	1:A:319:GLU:HG2	2.13	0.46
1:L:408:VAL:HA	1:L:411:ILE:HG22	1.97	0.46
1:L:200:GLN:OE1	1:L:221:GLU:HB3	2.15	0.46
1:A:303:ILE:HG23	1:A:338:PHE:HB3	1.98	0.46
3:G:133:LYS:HD2	3:G:164:VAL:HG21	1.96	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:113:ILE:HD13	1:L:326:LEU:CD2	2.45	0.46
1:A:182:ASN:ND2	8:A:569:HOH:O	2.48	0.46
1:A:408:VAL:O	1:A:411:ILE:HG22	2.15	0.46
2:B:85:LYS:HE3	2:B:224:LEU:HD13	1.99	0.46
1:L:385:ASP:O	1:L:389:VAL:HG23	2.15	0.46
1:A:23:TRP:HB3	1:A:28:PHE:HB2	1.97	0.45
1:A:151:PRO:HG3	8:A:569:HOH:O	2.16	0.45
2:E:149:SER:O	2:E:150:ASN:HB3	2.16	0.45
1:L:170:GLU:HB2	1:L:200:GLN:NE2	2.32	0.45
1:L:316:VAL:O	1:L:319:GLU:HG2	2.16	0.45
2:E:163:GLU:OE1	2:E:163:GLU:N	2.47	0.45
2:B:92:ILE:O	2:B:96:GLU:HG3	2.17	0.45
3:M:150:LYS:H	3:M:150:LYS:HD2	1.82	0.45
2:E:193:ARG:HG3	2:E:197:GLN:HB2	1.99	0.45
2:E:45:ASP:CG	2:E:46:GLY:N	2.70	0.45
2:B:47:PHE:HD1	2:B:47:PHE:H	1.64	0.45
2:E:204:HIS:HD2	8:E:227:HOH:O	1.99	0.45
1:L:229:GLU:HB3	1:L:230:PRO:HD3	1.99	0.45
1:L:534:GLN:HA	1:L:538:TRP:HB2	2.00	0.44
2:B:163:GLU:HG3	8:B:1070:HOH:O	2.17	0.44
1:A:331:ALA:HA	1:A:359:PHE:HB2	1.98	0.44
1:L:429:PRO:HA	1:L:430:PRO:HD3	1.84	0.44
3:M:55:THR:HG21	3:M:64:PHE:CZ	2.53	0.44
3:G:57:THR:OG1	3:G:59:LYS:HG2	2.18	0.44
1:A:202:SER:HB3	8:A:562:HOH:O	2.18	0.43
1:L:548:LEU:HD22	1:L:548:LEU:H	1.82	0.43
2:B:113:VAL:HG12	8:B:289:HOH:O	2.17	0.43
1:A:223:ILE:HG22	1:A:242:LYS:HD2	2.01	0.43
1:L:300:VAL:HG12	1:L:301:SER:N	2.34	0.43
1:L:549:ASP:OD2	1:L:551:ASN:HA	2.18	0.43
2:B:201:ALA:O	2:B:205:ILE:HG13	2.18	0.43
1:L:548:LEU:N	1:L:548:LEU:HD22	2.34	0.43
2:E:178:ARG:HA	2:E:178:ARG:HD3	1.82	0.43
1:A:62:ILE:HG23	1:A:286:ILE:HD11	2.01	0.42
1:L:160:GLU:HG3	1:L:408:VAL:HG13	2.02	0.42
2:B:157:ALA:N	2:B:158:PRO:CD	2.82	0.42
3:G:147:TYR:O	3:G:148:GLN:HB2	2.19	0.42
1:L:262:SER:HB2	1:L:276:MET:HB3	2.01	0.42
8:L:783:HOH:O	2:E:190:GLN:HG3	2.19	0.42
1:L:462:GLU:O	1:L:466:LYS:HB2	2.19	0.42
1:A:300:VAL:HG12	1:A:301:SER:H	1.85	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:1601:B12:H531	7:B:1601:B12:C55	2.45	0.42
1:A:310:PRO:HG2	1:L:542:LYS:HD2	2.01	0.42
1:A:305:VAL:HB	1:A:306:PRO:CD	2.50	0.42
1:A:365:SER:HB2	1:A:377:SER:HB3	2.02	0.42
3:M:134:GLU:H	3:M:134:GLU:CD	2.24	0.42
3:G:36:ARG:HD2	3:G:88:LEU:CD2	2.45	0.41
1:L:333:SER:HA	1:L:362:SER:OG	2.20	0.41
2:E:113:VAL:CG1	7:E:2601:B12:HM61	2.51	0.41
3:M:49:HIS:N	3:M:50:PRO:HD3	2.36	0.41
1:L:335:ASP:N	1:L:335:ASP:OD1	2.54	0.41
1:A:135:ARG:HD2	1:A:137:THR:O	2.20	0.41
7:B:1601:B12:C2B	7:B:1601:B12:H492	2.50	0.41
1:A:335:ASP:N	1:A:335:ASP:OD1	2.52	0.41
1:A:220:ALA:HB1	1:A:223:ILE:HD11	2.02	0.41
1:L:548:LEU:N	1:L:548:LEU:HD13	2.35	0.41
7:E:2601:B12:C2B	7:E:2601:B12:H492	2.50	0.41
1:L:331:ALA:HB1	1:L:362:SER:HB3	2.03	0.41
2:B:204:HIS:HD2	8:B:241:HOH:O	2.04	0.41
1:L:302:CYS:O	1:L:306:PRO:HD2	2.21	0.41
2:B:153:LEU:HG	2:B:155:PRO:HD3	2.03	0.40
1:L:209:LEU:O	1:L:213:MET:HG3	2.21	0.40
1:L:428:LEU:O	1:L:429:PRO:C	2.58	0.40
1:L:153:GLN:OE1	1:L:445:SER:HA	2.21	0.40
1:A:162:ALA:HB3	1:A:193:GLY:HA3	2.02	0.40
2:B:100:ILE:HD11	2:B:177:LYS:HG3	2.04	0.40
1:L:266:VAL:HG23	8:L:585:HOH:O	2.20	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/554 (99%)	526 (96%)	22 (4%)	2 (0%)	34	32
1	L	549/554 (99%)	516 (94%)	31 (6%)	2 (0%)	34	32
2	B	177/224 (79%)	173 (98%)	4 (2%)	0	100	100
2	E	178/224 (80%)	175 (98%)	3 (2%)	0	100	100
3	G	137/173 (79%)	135 (98%)	2 (2%)	0	100	100
3	M	136/173 (79%)	132 (97%)	4 (3%)	0	100	100
All	All	1727/1902 (91%)	1657 (96%)	66 (4%)	4 (0%)	47	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	300	VAL
1	A	363	GLY
1	L	300	VAL
1	L	363	GLY

### 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	451/453 (100%)	443 (98%)	8 (2%)	59	65
1	L	450/453 (99%)	439 (98%)	11 (2%)	49	53
2	B	147/183 (80%)	146 (99%)	1 (1%)	84	88
2	E	148/183 (81%)	148 (100%)	0	100	100
3	G	118/141 (84%)	116 (98%)	2 (2%)	60	67
3	M	117/141 (83%)	115 (98%)	2 (2%)	60	67
All	All	1431/1554 (92%)	1407 (98%)	24 (2%)	60	67

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	PHE

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	A	178	TYR
1	A	182	ASN
1	A	232	PHE
1	A	348	ARG
1	A	501	TYR
1	A	548	LEU
1	A	551	ASN
2	B	167	GLN
3	G	58	ASN
3	G	131	SER
1	L	1	MET
1	L	2	ARG
1	L	25	GLU
1	L	178	TYR
1	L	182	ASN
1	L	232	PHE
1	L	348	ARG
1	L	350	LEU
1	L	501	TYR
1	L	534	GLN
1	L	548	LEU
3	M	58	ASN
3	M	150	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	182	ASN
1	A	465	ASN
1	A	479	GLN
1	A	551	ASN
2	B	142	GLN
2	B	143	GLN
2	B	204	HIS
3	G	58	ASN
1	L	35	ASN
1	L	45	ASN
1	L	182	ASN
1	L	465	ASN
1	L	467	ASN
1	L	551	ASN
2	E	143	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	E	167	GLN
2	E	204	HIS
2	E	218	GLN
3	M	58	ASN

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
6	PO4	L	1402	-	4,4,4	1.67	0	6,6,6	0.43	0
6	PO4	M	1404	-	4,4,4	1.66	0	6,6,6	0.44	0
4	GOL	L	2602	5	5,5,5	0.22	0	5,5,5	0.16	0
4	GOL	A	1602	5	5,5,5	0.23	0	5,5,5	0.16	0
7	B12	E	2601	-	80,101,101	1.20	7 (8%)	101,166,166	1.26	11 (10%)
7	B12	B	1601	-	80,101,101	1.23	6 (7%)	101,166,166	1.23	10 (9%)
6	PO4	E	1405	-	4,4,4	1.55	0	6,6,6	0.42	0
6	PO4	B	1406	-	4,4,4	1.59	0	6,6,6	0.43	0
6	PO4	A	1401	-	4,4,4	1.66	0	6,6,6	0.42	0
6	PO4	G	1403	-	4,4,4	1.66	0	6,6,6	0.44	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
7	B12	E	2601	-	-	7/51/223/223	0/3/11/11
7	B12	B	1601	-	-	8/51/223/223	0/3/11/11
4	GOL	L	2602	5	-	0/4/4/4	-
4	GOL	A	1602	5	-	0/4/4/4	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	E	2601	B12	C5M-C5B	-5.56	1.39	1.51
7	B	1601	B12	C5M-C5B	-5.52	1.40	1.51
7	B	1601	B12	C11-C10	-3.61	1.34	1.40
7	E	2601	B12	C11-C10	-3.36	1.35	1.40
7	E	2601	B12	C17-C18	2.98	1.58	1.54
7	B	1601	B12	C17-C18	2.91	1.58	1.54
7	B	1601	B12	C1-C2	2.75	1.65	1.58
7	E	2601	B12	O58-C57	2.63	1.28	1.23
7	B	1601	B12	O58-C57	2.57	1.28	1.23
7	B	1601	B12	C48-C49	2.56	1.60	1.52
7	E	2601	B12	C48-C49	2.54	1.60	1.52
7	E	2601	B12	C1-C2	2.42	1.64	1.58
7	E	2601	B12	C16-C15	2.04	1.48	1.41

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	E	2601	B12	C26-C2-C1	3.78	115.90	110.02
7	B	1601	B12	C26-C2-C1	3.59	115.61	110.02
7	E	2601	B12	C55-C17-C16	3.53	121.68	109.92
7	B	1601	B12	C55-C17-C16	3.48	121.50	109.92
7	B	1601	B12	C16-C15-C14	-3.25	119.20	124.27
7	E	2601	B12	C16-C15-C14	-3.20	119.28	124.27
7	E	2601	B12	C56-C55-C17	2.94	121.19	115.50
7	B	1601	B12	C56-C55-C17	2.85	121.02	115.50
7	B	1601	B12	O58-C57-C56	-2.57	117.31	122.02
7	E	2601	B12	O58-C57-C56	-2.48	117.49	122.02
7	E	2601	B12	C47-C12-C46	-2.43	104.63	109.73
7	B	1601	B12	C4B-C9B-C8B	-2.41	118.64	121.10
7	B	1601	B12	C7-C37-C38	2.40	121.33	114.20

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1601	B12	C47-C12-C46	-2.37	104.76	109.73
7	B	1601	B12	C8-C9-N22	2.30	113.99	111.12
7	E	2601	B12	C7-C37-C38	2.29	121.02	114.20
7	E	2601	B12	C8-C9-N22	2.28	113.97	111.12
7	E	2601	B12	C19-C1-N21	2.25	104.47	102.16
7	E	2601	B12	C4B-C9B-C8B	-2.21	118.83	121.10
7	E	2601	B12	C20-C1-C19	-2.21	107.23	109.36
7	B	1601	B12	C3-C4-C5	-2.01	124.40	131.68

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	E	2601	B12	C12-C13-C48-C49
7	B	1601	B12	C12-C13-C48-C49
7	E	2601	B12	C1-C2-C26-C27
7	E	2601	B12	C25-C2-C26-C27
7	B	1601	B12	C1-C2-C26-C27
7	B	1601	B12	C25-C2-C26-C27
7	E	2601	B12	C4-C3-C30-C31
7	B	1601	B12	C4-C3-C30-C31
7	E	2601	B12	C17-C18-C60-C61
7	B	1601	B12	C17-C18-C60-C61
7	E	2601	B12	C2-C3-C30-C31
7	E	2601	B12	C19-C18-C60-C61
7	B	1601	B12	C19-C18-C60-C61
7	B	1601	B12	C2-C3-C30-C31
7	B	1601	B12	C13-C48-C49-C50

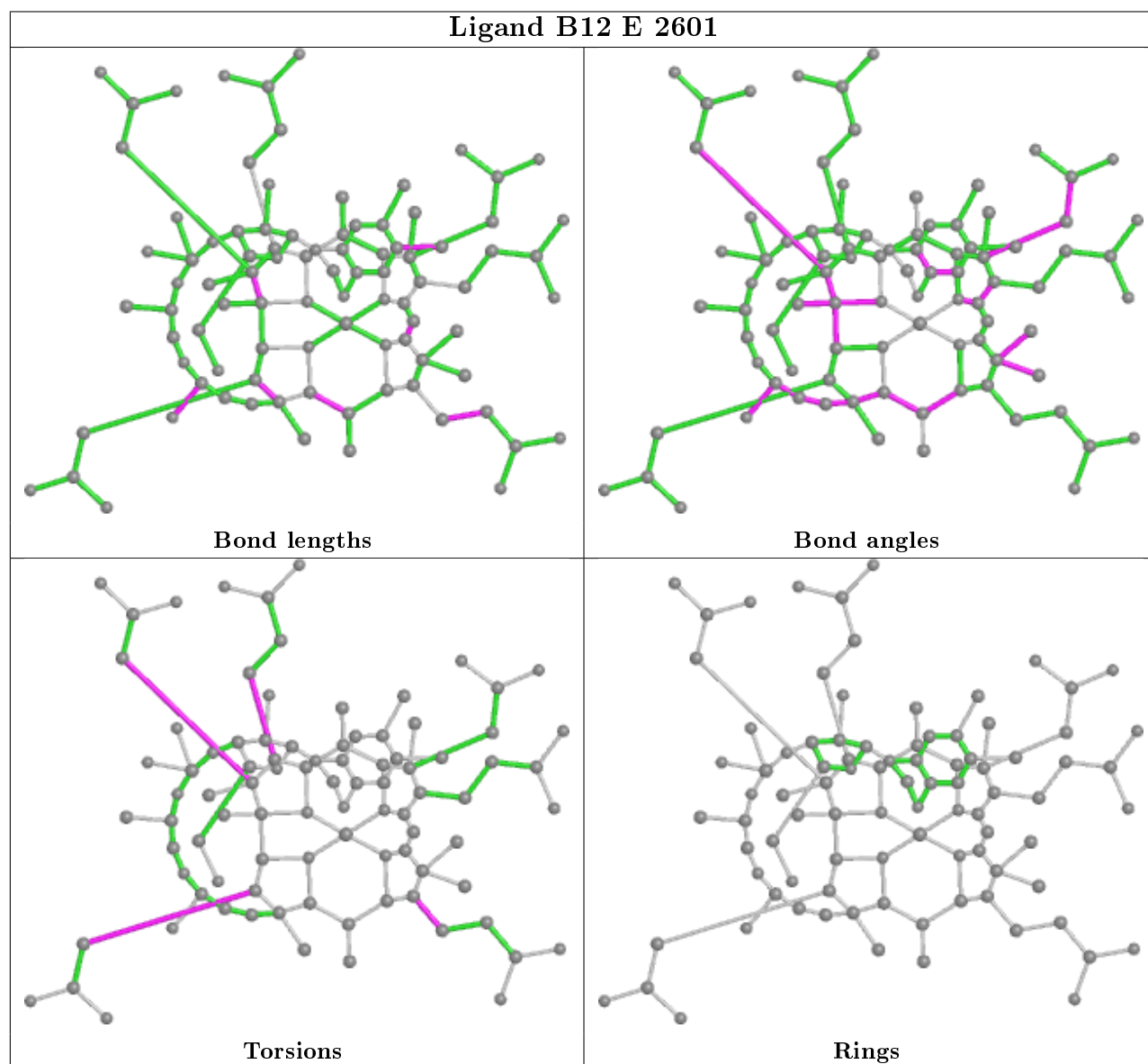
There are no ring outliers.

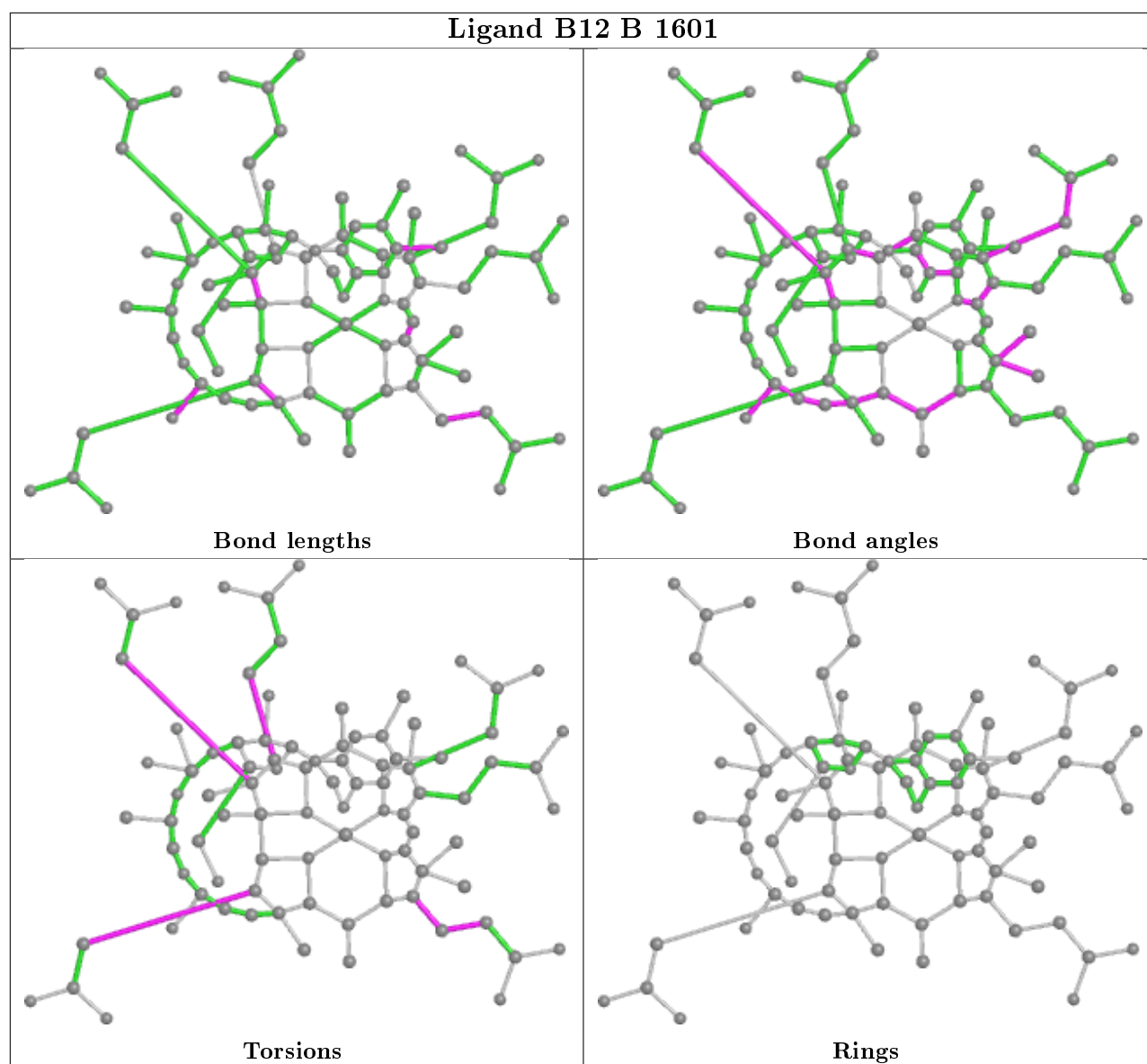
2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	E	2601	B12	5	0
7	B	1601	B12	6	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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	552/554 (99%)	-0.34	4 (0%) 87 89	17, 24, 36, 83	0
1	L	551/554 (99%)	-0.26	8 (1%) 73 77	17, 24, 37, 69	0
2	B	179/224 (79%)	-0.48	2 (1%) 80 84	18, 27, 38, 63	0
2	E	180/224 (80%)	-0.52	3 (1%) 70 74	20, 26, 39, 84	0
3	G	139/173 (80%)	-0.19	2 (1%) 75 78	21, 31, 47, 85	0
3	M	138/173 (79%)	-0.09	2 (1%) 75 78	23, 33, 55, 83	0
All	All	1739/1902 (91%)	-0.32	21 (1%) 79 82	17, 26, 42, 85	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	35	SER	7.9
2	B	224	LEU	5.9
2	E	45	ASP	4.5
1	A	1	MET	4.0
1	A	552	GLU	3.6
2	E	224	LEU	3.5
1	L	551	ASN	3.4
3	M	36	ARG	3.3
2	E	46	GLY	3.0
1	L	222	THR	2.7
1	L	223	ILE	2.7
2	B	46	GLY	2.6
1	L	548	LEU	2.5
1	L	259	THR	2.4
1	L	295	LEU	2.3
3	G	36	ARG	2.3
1	L	1	MET	2.2
1	A	202	SER	2.2
1	A	223	ILE	2.2

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
3	M	77	GLN	2.0
1	L	173	VAL	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 carbohydrates 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

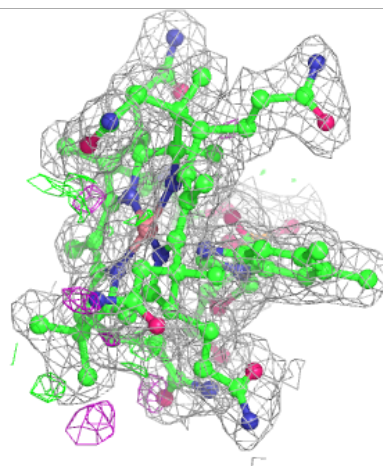
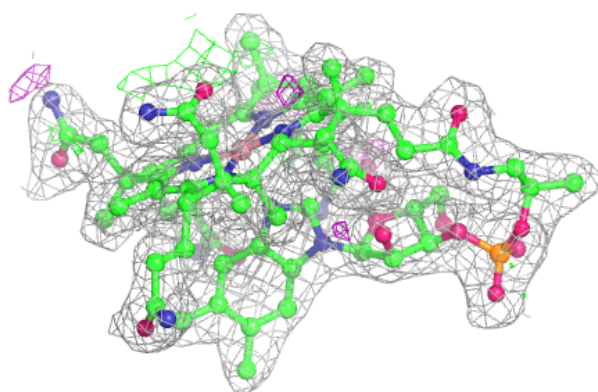
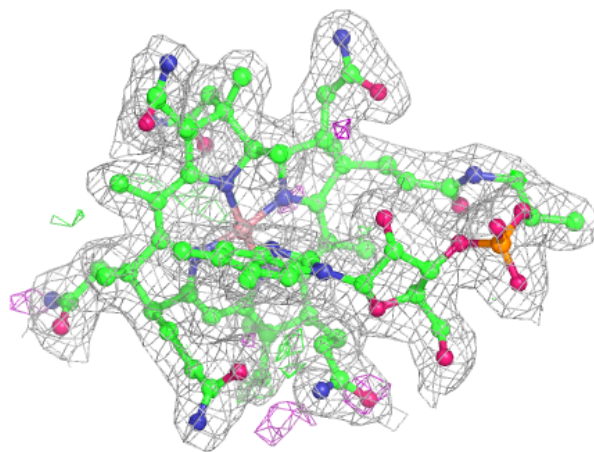
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	PO4	E	1405	5/5	0.91	0.30	70,71,72,72	0
6	PO4	B	1406	5/5	0.93	0.31	70,71,72,72	0
6	PO4	M	1404	5/5	0.94	0.13	61,61,62,63	0
4	GOL	L	2602	6/6	0.95	0.21	20,24,28,34	0
6	PO4	A	1401	5/5	0.95	0.14	74,74,75,75	0
6	PO4	L	1402	5/5	0.95	0.23	78,78,78,78	0
7	B12	B	1601	91/91	0.97	0.14	16,20,24,25	0
4	GOL	A	1602	6/6	0.97	0.17	15,20,26,33	0
7	B12	E	2601	91/91	0.97	0.14	15,20,24,26	0
6	PO4	G	1403	5/5	0.97	0.12	61,61,61,62	0
5	CA	L	2603	1/1	0.99	0.11	21,21,21,21	0
5	CA	A	1603	1/1	1.00	0.10	20,20,20,20	0

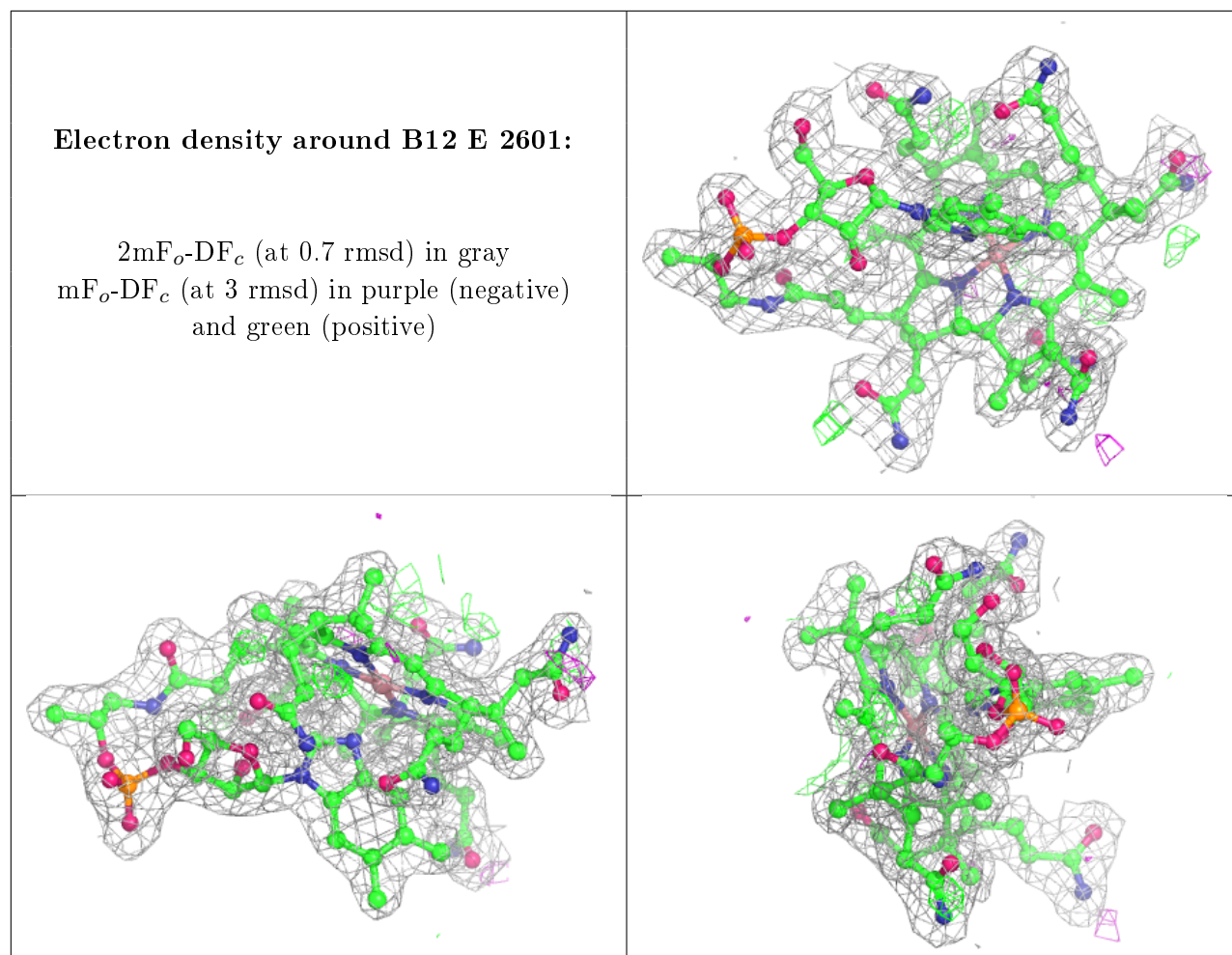
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around B12 B 1601:**

$2mF_o-DF_c$  (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.