



# Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 08:08 am BST

PDB ID : 2E93  
Title : S. cerevisiae geranylgeranyl pyrophosphate synthase in complex with BPH-629  
Authors : Guo, R.T.; Cao, R.; Ko, T.P.; Chen, C.K.-M.; Jeng, W.Y.; Chang, T.H.;  
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Deposited on : 2007-01-24  
Resolution : 2.12 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.

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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

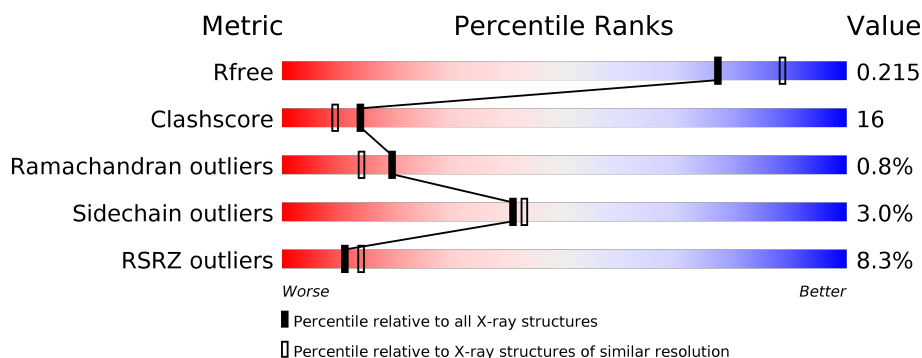
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.12 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	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)

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	340	<div> <div>5%</div> <div> <div></div> <div>70%</div> <div>16%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	340	<div> <div>10%</div> <div> <div></div> <div>68%</div> <div>21%</div> <div>•</div> <div>8%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5652 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.

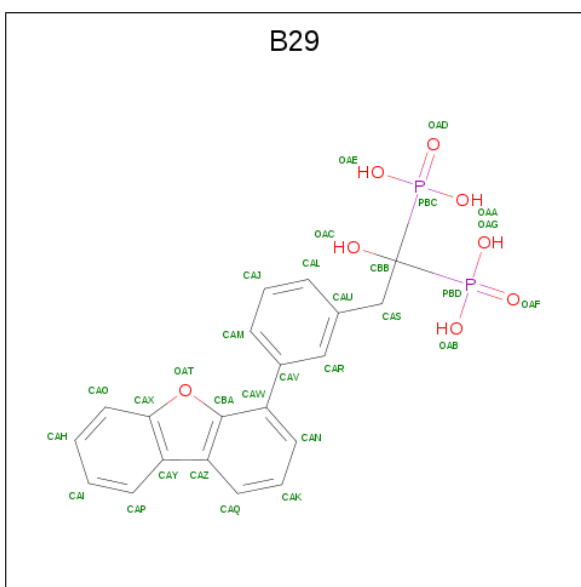
- Molecule 1 is a protein called Geranylgeranyl pyrophosphate synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2442	1569	414	450	9			
1	B	314	Total	C	N	O	S	0	0	0
			2571	1652	432	477	10			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP Q12051
A	2	THR	-	CLONING ARTIFACT	UNP Q12051
A	3	LYS	-	CLONING ARTIFACT	UNP Q12051
A	4	ASN	-	CLONING ARTIFACT	UNP Q12051
A	5	LYS	-	CLONING ARTIFACT	UNP Q12051
B	1	MET	-	CLONING ARTIFACT	UNP Q12051
B	2	THR	-	CLONING ARTIFACT	UNP Q12051
B	3	LYS	-	CLONING ARTIFACT	UNP Q12051
B	4	ASN	-	CLONING ARTIFACT	UNP Q12051
B	5	LYS	-	CLONING ARTIFACT	UNP Q12051

- Molecule 2 is [2-(3-DIBENZOFURAN-4-YL-PHENYL)-1-HYDROXY-1-PHOSPHONO-ETHYL]-PHOSPHONIC ACID (three-letter code: B29) (formula: C<sub>20</sub>H<sub>18</sub>O<sub>8</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 30	C 20	O 8	P 2	0	0
2	A	1	Total 30	C 20	O 8	P 2	0	0
2	B	1	Total 30	C 20	O 8	P 2	0	0
2	B	1	Total 30	C 20	O 8	P 2	0	0

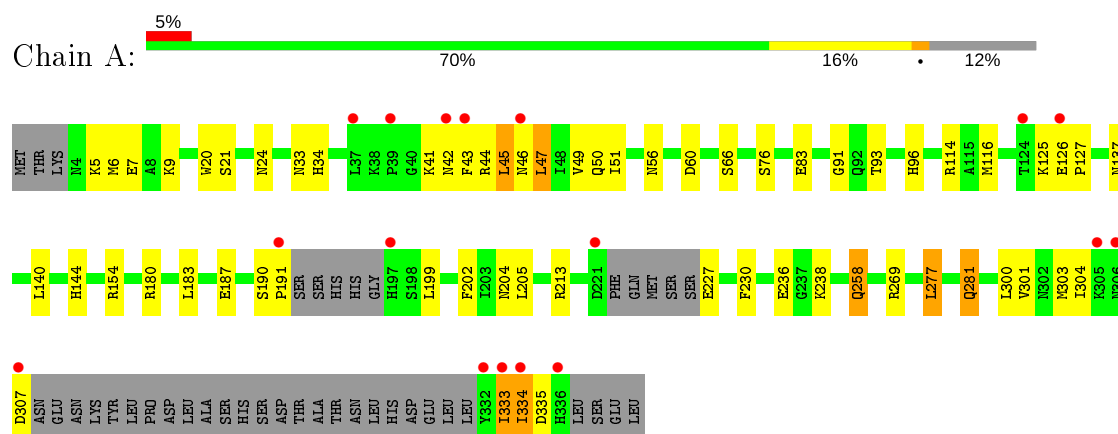
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	280	Total O 280 280	0	0
3	B	239	Total O 239 239	0	0

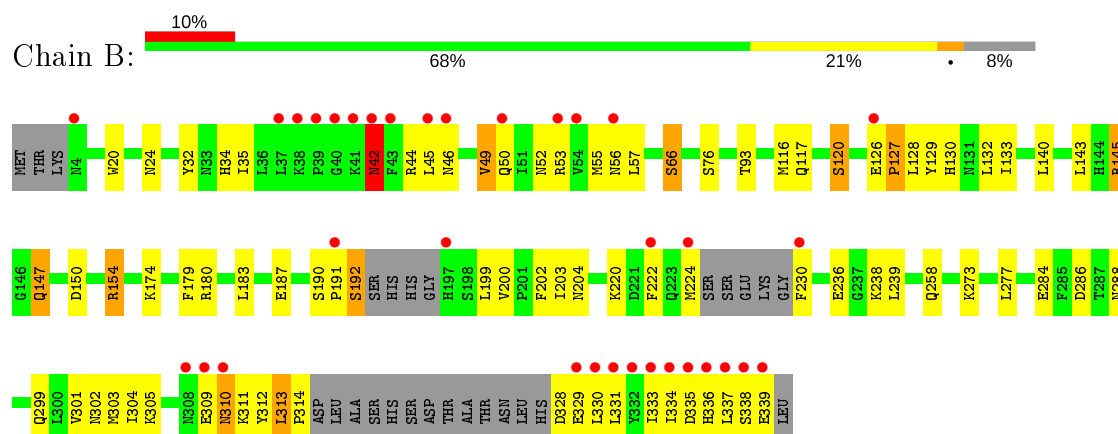
### 3 Residue-property plots [i](#)

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: Geranylgeranyl pyrophosphate synthetase



- Molecule 1: Geranylgeranyl pyrophosphate synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.79Å 115.82Å 130.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.96 – 2.12 28.96 – 2.12	Depositor EDS
% Data completeness (in resolution range)	93.6 (28.96-2.12) 94.1 (28.96-2.12)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.11 (at 2.12Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.167 , 0.217 0.165 , 0.215	Depositor DCC
$R_{free}$ test set	1873 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.2	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 63.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5652	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% 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: B29

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	0.90	1/2485 (0.0%)	0.88	4/3362 (0.1%)
1	B	0.87	0/2617	0.83	3/3542 (0.1%)
All	All	0.88	1/5102 (0.0%)	0.86	7/6904 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	GLU	CB-CG	5.06	1.61	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ARG	NE-CZ-NH2	-10.00	115.30	120.30
1	A	60	ASP	CB-CG-OD1	8.07	125.57	118.30
1	A	269	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	B	145	ARG	NE-CZ-NH1	-5.76	117.42	120.30
1	B	150	ASP	CB-CG-OD1	5.59	123.33	118.30
1	B	154	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	A	277	LEU	CA-CB-CG	-5.14	103.48	115.30

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	2442	0	2494	73	0
1	B	2571	0	2615	96	0
2	A	60	0	28	6	0
2	B	60	0	28	9	0
3	A	280	0	0	12	0
3	B	239	0	0	8	0
All	All	5652	0	5165	163	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 16.

All (163) 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:B:304:ILE:HG21	1:B:334:ILE:HD13	1.28	1.14
1:B:126:GLU:HG3	3:B:1089:HOH:O	1.69	0.92
1:B:258:GLN:HG2	1:B:286:ASP:OD2	1.69	0.91
1:B:52:ASN:HD21	1:B:57:LEU:H	1.12	0.89
1:A:258:GLN:H	1:A:258:GLN:HE21	1.16	0.89
1:A:205:LEU:HD23	1:A:304:ILE:HD11	1.57	0.85
1:B:301:VAL:HG13	1:B:334:ILE:HG22	1.59	0.85
1:A:205:LEU:HD23	1:A:304:ILE:CD1	2.11	0.80
1:B:127:PRO:HG2	3:B:1090:HOH:O	1.85	0.77
1:A:281:GLN:HA	1:A:281:GLN:HE21	1.52	0.75
1:A:45:LEU:HD11	1:A:66:SER:HB3	1.69	0.75
1:A:43:PHE:O	1:A:46:ASN:HB3	1.86	0.74
1:B:313:LEU:HD11	1:B:330:LEU:HD23	1.68	0.73
1:B:49:VAL:HG12	1:B:50:GLN:HE21	1.51	0.73
1:B:304:ILE:HG22	1:B:331:LEU:HD13	1.72	0.72
1:B:55:MET:CE	1:B:187:GLU:HA	2.21	0.69
1:B:331:LEU:HA	1:B:334:ILE:HD12	1.74	0.69
1:B:45:LEU:HD13	1:B:66:SER:OG	1.93	0.69
1:A:304:ILE:CB	1:A:334:ILE:HD12	2.23	0.69
1:A:304:ILE:HB	1:A:334:ILE:HD12	1.75	0.69
1:B:180:ARG:HH11	1:B:204:ASN:HD21	1.40	0.68
1:A:44:ARG:HG3	2:A:902:B29:CAX	2.23	0.68
1:A:116:MET:CE	1:B:116:MET:SD	2.82	0.68
1:A:334:ILE:O	1:A:334:ILE:HG12	1.92	0.68
1:B:52:ASN:ND2	1:B:57:LEU:H	1.92	0.65
1:A:227:GLU:N	3:A:1121:HOH:O	2.29	0.65
1:A:116:MET:HE1	1:B:116:MET:SD	2.38	0.64
1:A:230:PHE:HA	3:A:1067:HOH:O	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:MET:HE2	1:B:187:GLU:HA	1.80	0.64
1:B:55:MET:HE3	1:B:190:SER:HB2	1.79	0.64
1:B:180:ARG:HH11	1:B:204:ASN:ND2	1.96	0.63
1:B:230:PHE:HZ	3:B:1023:HOH:O	1.81	0.63
1:A:7:GLU:HG3	1:B:303:MET:HG2	1.80	0.63
3:A:995:HOH:O	1:B:130:HIS:HE1	1.79	0.62
1:A:91:GLY:HA2	1:A:227:GLU:HB3	1.81	0.62
1:B:313:LEU:CD1	1:B:330:LEU:HD23	2.29	0.62
1:A:6:MET:SD	1:B:299:GLN:NE2	2.73	0.62
1:A:43:PHE:C	1:A:46:ASN:HB3	2.20	0.61
1:A:304:ILE:HB	1:A:334:ILE:CD1	2.31	0.60
1:B:335:ASP:O	1:B:338:SER:HB2	2.01	0.60
1:B:220:LYS:O	1:B:224:MET:HG2	2.01	0.60
1:B:50:GLN:OE1	1:B:329:GLU:HG2	2.02	0.59
1:B:330:LEU:HA	1:B:333:ILE:HD12	1.83	0.59
1:B:53:ARG:HD2	3:B:1108:HOH:O	2.03	0.58
1:B:147:GLN:HA	1:B:147:GLN:HE21	1.69	0.57
1:A:238:LYS:NZ	1:A:238:LYS:HB3	2.18	0.57
1:A:126:GLU:HB2	1:A:127:PRO:HD3	1.85	0.57
1:A:140:LEU:O	1:A:144:HIS:HD2	1.88	0.56
1:A:44:ARG:HD2	2:A:902:B29:CAZ	2.35	0.56
1:B:126:GLU:N	1:B:127:PRO:HD2	2.19	0.56
1:B:55:MET:HE3	1:B:190:SER:CB	2.35	0.56
1:B:309:GLU:O	1:B:310:ASN:HB2	2.05	0.56
1:B:273:LYS:NZ	1:B:277:LEU:HD11	2.20	0.56
1:A:333:ILE:O	1:A:335:ASP:N	2.38	0.56
1:B:55:MET:HE1	1:B:187:GLU:HA	1.87	0.56
1:A:213:ARG:NH1	3:A:1160:HOH:O	2.35	0.56
1:B:339:GLU:HA	1:B:339:GLU:OE2	2.05	0.56
1:A:6:MET:O	1:A:9:LYS:HG2	2.05	0.56
1:A:21:SER:H	1:A:24:ASN:HD22	1.54	0.56
1:A:258:GLN:HE21	1:A:258:GLN:N	1.96	0.55
1:A:281:GLN:HA	1:A:281:GLN:NE2	2.18	0.55
1:B:273:LYS:HZ1	1:B:277:LEU:HD11	1.72	0.55
1:A:180:ARG:HH11	1:A:204:ASN:HD21	1.53	0.54
1:B:120:SER:HA	1:B:129:TYR:HE1	1.73	0.54
1:B:179:PHE:HB3	1:B:203:ILE:HG13	1.89	0.53
1:A:180:ARG:HH11	1:A:204:ASN:ND2	2.06	0.53
1:A:227:GLU:N	3:A:1142:HOH:O	2.40	0.53
1:B:183:LEU:HD21	1:B:199:LEU:C	2.28	0.53
1:A:281:GLN:HE21	1:A:281:GLN:CA	2.21	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:ASN:O	1:A:50:GLN:HG2	2.09	0.53
1:A:116:MET:HE2	1:B:116:MET:SD	2.49	0.52
1:B:55:MET:HE1	1:B:187:GLU:HG2	1.91	0.52
1:B:55:MET:CE	1:B:190:SER:HB2	2.39	0.52
1:A:45:LEU:O	1:A:49:VAL:HG23	2.09	0.52
2:B:903:B29:OAT	2:B:903:B29:HAM	2.08	0.52
1:B:143:LEU:HD23	1:B:174:LYS:O	2.10	0.52
1:B:34:HIS:HE1	1:B:93:THR:O	1.93	0.52
1:B:304:ILE:HD13	1:B:334:ILE:HG12	1.91	0.51
1:A:202:PHE:CE1	1:A:304:ILE:HD12	2.45	0.51
1:A:180:ARG:HD3	1:A:204:ASN:HD21	1.76	0.51
1:A:83:GLU:O	1:A:154:ARG:NH2	2.43	0.51
1:B:202:PHE:CE1	1:B:334:ILE:HD11	2.45	0.51
1:A:114:ARG:NH1	3:A:931:HOH:O	2.43	0.51
1:A:43:PHE:HA	1:A:46:ASN:HB2	1.93	0.51
1:A:5:LYS:HD2	3:A:907:HOH:O	2.12	0.50
1:B:56:ASN:ND2	1:B:192:SER:OG	2.41	0.50
1:B:304:ILE:HG23	1:B:313:LEU:HD22	1.93	0.50
1:A:154:ARG:NH1	3:A:977:HOH:O	2.44	0.49
1:B:284:GLU:HG2	3:B:1058:HOH:O	2.11	0.49
1:B:191:PRO:HG2	1:B:192:SER:H	1.77	0.49
1:A:20:TRP:CH2	1:A:24:ASN:HB3	2.48	0.49
1:B:222:PHE:HD1	1:B:230:PHE:N	2.09	0.49
1:B:76:SER:HB2	2:B:904:B29:CAY	2.42	0.49
1:B:20:TRP:CH2	1:B:24:ASN:HB3	2.48	0.49
1:A:334:ILE:CG1	1:A:334:ILE:O	2.61	0.48
1:A:96:HIS:HD2	3:A:966:HOH:O	1.95	0.48
1:B:44:ARG:HH11	2:B:903:B29:HAJ	1.78	0.48
1:B:76:SER:HB2	2:B:904:B29:CAZ	2.43	0.48
1:A:304:ILE:HG13	1:A:334:ILE:HD12	1.96	0.48
1:B:202:PHE:CD1	1:B:313:LEU:HD13	2.49	0.48
1:A:33:ASN:ND2	3:A:1113:HOH:O	2.36	0.48
1:A:47:LEU:HG	1:A:51:ILE:HD12	1.96	0.48
1:B:147:GLN:HE22	1:B:174:LYS:NZ	2.11	0.48
1:B:329:GLU:O	1:B:333:ILE:HG13	2.13	0.48
1:B:76:SER:HB2	2:B:904:B29:CAX	2.44	0.48
1:B:273:LYS:NZ	1:B:277:LEU:CD1	2.77	0.47
1:B:202:PHE:CZ	1:B:334:ILE:HD11	2.48	0.47
1:A:137:ASN:HD21	1:B:117:GLN:N	2.12	0.47
1:A:34:HIS:HE1	1:A:93:THR:O	1.97	0.47
1:B:55:MET:O	1:B:191:PRO:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LYS:HD2	3:A:1146:HOH:O	2.15	0.46
1:A:301:VAL:HG13	1:A:334:ILE:HD11	1.97	0.46
1:A:76:SER:HB2	2:A:901:B29:CAZ	2.45	0.46
1:A:42:ASN:C	1:A:44:ARG:H	2.18	0.46
1:B:313:LEU:HD12	1:B:314:PRO:CD	2.45	0.46
1:A:300:LEU:O	1:A:303:MET:HB2	2.15	0.46
1:A:43:PHE:HA	1:A:46:ASN:CB	2.46	0.46
1:A:277:LEU:HA	1:A:277:LEU:HD12	1.84	0.45
1:A:76:SER:HB2	2:A:901:B29:CAY	2.46	0.45
1:B:313:LEU:HD12	1:B:314:PRO:HD2	1.98	0.45
1:B:76:SER:HB2	2:B:904:B29:CBA	2.47	0.45
1:B:154:ARG:HD2	1:B:236:GLU:O	2.17	0.45
1:A:44:ARG:HD3	1:A:44:ARG:HA	1.45	0.44
1:A:304:ILE:CG2	1:A:334:ILE:HD12	2.47	0.44
2:A:901:B29:OAE	2:A:902:B29:OAE	2.35	0.44
1:B:32:TYR:O	1:B:35:ILE:HG22	2.18	0.44
1:B:313:LEU:HA	1:B:313:LEU:HD12	1.90	0.44
1:A:20:TRP:CZ3	1:B:145:ARG:HB3	2.53	0.44
1:A:187:GLU:OE2	3:A:1049:HOH:O	2.21	0.43
1:B:183:LEU:HD23	1:B:200:VAL:HA	2.00	0.43
1:B:301:VAL:HG13	1:B:334:ILE:CG2	2.39	0.43
1:B:309:GLU:O	1:B:310:ASN:CB	2.66	0.43
1:A:304:ILE:HG21	1:A:334:ILE:HD12	1.99	0.43
1:B:202:PHE:CZ	1:B:330:LEU:HG	2.52	0.43
1:B:310:ASN:HB3	1:B:313:LEU:O	2.18	0.43
1:A:44:ARG:HG3	2:A:902:B29:OAT	2.18	0.43
1:B:239:LEU:HA	1:B:239:LEU:HD13	1.81	0.43
1:B:302:ASN:HA	1:B:305:LYS:HB2	2.01	0.43
1:A:126:GLU:HB3	1:B:126:GLU:HB3	2.01	0.43
1:B:174:LYS:HE2	2:B:904:B29:OAG	2.18	0.43
1:B:46:ASN:O	1:B:50:GLN:HG2	2.19	0.42
1:B:44:ARG:HH11	2:B:903:B29:CAJ	2.32	0.42
2:B:903:B29:OAE	2:B:904:B29:OAA	2.37	0.42
1:A:304:ILE:CG1	1:A:334:ILE:HD12	2.49	0.42
1:B:238:LYS:CE	3:B:1014:HOH:O	2.67	0.42
1:B:238:LYS:HE3	3:B:1014:HOH:O	2.19	0.42
1:B:52:ASN:HD22	1:B:55:MET:HB2	1.83	0.42
1:A:47:LEU:O	1:A:51:ILE:HD12	2.20	0.42
1:B:128:LEU:O	1:B:132:LEU:HG	2.19	0.42
1:B:133:ILE:HA	1:B:133:ILE:HD13	1.80	0.42
1:A:190:SER:HA	1:A:191:PRO:HD3	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:GLU:N	1:B:127:PRO:CD	2.82	0.42
1:B:311:LYS:HE3	1:B:312:TYR:CZ	2.55	0.42
1:A:183:LEU:HD21	1:A:199:LEU:HB2	2.02	0.42
1:B:183:LEU:HD22	1:B:203:ILE:HB	2.01	0.42
1:B:130:HIS:O	1:B:130:HIS:HD2	2.03	0.41
1:B:333:ILE:O	1:B:337:LEU:HG	2.20	0.41
1:B:311:LYS:HE3	1:B:312:TYR:CE2	2.55	0.41
1:A:144:HIS:HE1	3:B:922:HOH:O	2.03	0.41
1:B:42:ASN:N	1:B:42:ASN:HD22	2.19	0.41
1:A:41:LYS:HA	1:A:41:LYS:HD3	1.89	0.40
1:B:180:ARG:HD3	1:B:204:ASN:HD21	1.86	0.40
1:A:137:ASN:CB	1:B:116:MET:HE1	2.52	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	291/340 (86%)	282 (97%)	7 (2%)	2 (1%)	22	17
1	B	306/340 (90%)	299 (98%)	4 (1%)	3 (1%)	15	10
All	All	597/680 (88%)	581 (97%)	11 (2%)	5 (1%)	19	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	334	ILE
1	B	42	ASN
1	B	310	ASN
1	B	127	PRO
1	A	333	ILE

### 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	279/317 (88%)	273 (98%)	6 (2%)	52	55
1	B	295/317 (93%)	284 (96%)	11 (4%)	34	34
All	All	574/634 (90%)	557 (97%)	17 (3%)	41	43

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	A	47	LEU
1	A	56	ASN
1	A	258	GLN
1	A	281	GLN
1	A	307	ASP
1	B	42	ASN
1	B	49	VAL
1	B	66	SER
1	B	120	SER
1	B	140	LEU
1	B	147	GLN
1	B	192	SER
1	B	288	ASN
1	B	313	LEU
1	B	328	ASP
1	B	336	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	34	HIS
1	A	61	GLN
1	A	67	GLN
1	A	85	ASN

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Mol	Chain	Res	Type
1	A	96	HIS
1	A	109	ASN
1	A	137	ASN
1	A	144	HIS
1	A	164	GLN
1	A	204	ASN
1	A	258	GLN
1	A	260	ASN
1	A	281	GLN
1	B	24	ASN
1	B	34	HIS
1	B	42	ASN
1	B	52	ASN
1	B	56	ASN
1	B	85	ASN
1	B	96	HIS
1	B	109	ASN
1	B	130	HIS
1	B	147	GLN
1	B	204	ASN
1	B	288	ASN
1	B	298	ASN
1	B	299	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	B29	A	901	-	29,33,33	2.70	10 (34%)	43,52,52	1.17	5 (11%)
2	B29	A	902	-	29,33,33	2.70	10 (34%)	43,52,52	1.51	12 (27%)
2	B29	B	904	-	29,33,33	2.69	10 (34%)	43,52,52	1.23	4 (9%)
2	B29	B	903	-	29,33,33	2.72	10 (34%)	43,52,52	1.18	3 (6%)

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	B29	A	901	-	-	3/27/27/27	0/4/4/4
2	B29	A	902	-	-	5/27/27/27	0/4/4/4
2	B29	B	904	-	-	1/27/27/27	0/4/4/4
2	B29	B	903	-	-	5/27/27/27	0/4/4/4

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	903	B29	PBC-OAD	7.04	1.61	1.50
2	A	902	B29	PBD-OAF	6.93	1.61	1.50
2	B	904	B29	PBC-OAD	6.86	1.61	1.50
2	B	904	B29	PBD-OAF	6.85	1.61	1.50
2	B	903	B29	PBD-OAF	6.82	1.61	1.50
2	A	902	B29	PBC-OAD	6.80	1.61	1.50
2	A	901	B29	PBD-OAF	6.71	1.61	1.50
2	A	901	B29	PBC-OAD	6.62	1.60	1.50
2	B	903	B29	CAW-CAV	-5.83	1.39	1.49
2	A	901	B29	CAW-CAV	-5.77	1.39	1.49
2	B	904	B29	CAW-CAV	-5.54	1.39	1.49
2	A	902	B29	CAW-CAV	-5.53	1.39	1.49
2	B	904	B29	PBD-OAG	3.89	1.61	1.54
2	A	901	B29	PBD-OAG	3.76	1.61	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	903	B29	PBC-OAE	3.69	1.61	1.54
2	B	903	B29	PBD-OAG	3.60	1.61	1.54
2	A	901	B29	PBD-OAB	-3.60	1.48	1.54
2	B	904	B29	PBD-OAB	-3.59	1.48	1.54
2	A	901	B29	PBC-OAE	3.58	1.61	1.54
2	A	902	B29	PBC-OAA	-3.51	1.48	1.54
2	A	902	B29	PBC-OAE	3.51	1.61	1.54
2	B	904	B29	PBC-OAA	-3.48	1.48	1.54
2	B	903	B29	PBC-OAA	-3.45	1.48	1.54
2	A	902	B29	PBD-OAB	-3.43	1.48	1.54
2	B	904	B29	PBC-OAE	3.40	1.61	1.54
2	A	902	B29	PBC-CBB	3.40	1.87	1.85
2	B	903	B29	PBD-OAB	-3.38	1.48	1.54
2	A	901	B29	PBC-OAA	-3.38	1.48	1.54
2	A	902	B29	PBD-OAG	3.23	1.60	1.54
2	A	901	B29	PBD-CBB	2.90	1.87	1.85
2	B	904	B29	PBD-CBB	2.64	1.87	1.85
2	A	901	B29	PBC-CBB	2.32	1.86	1.85
2	A	902	B29	PBD-CBB	2.32	1.86	1.85
2	B	903	B29	PBC-CBB	2.31	1.86	1.85
2	B	903	B29	PBD-CBB	2.30	1.86	1.85
2	A	901	B29	CAZ-CAY	-2.17	1.39	1.45
2	A	902	B29	CAZ-CAY	-2.14	1.39	1.45
2	B	903	B29	CAZ-CAY	-2.09	1.39	1.45
2	B	904	B29	CAZ-CAY	-2.06	1.39	1.45
2	B	904	B29	PBC-CBB	2.06	1.86	1.85

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	904	B29	OAG-PBD-OAF	-3.85	104.41	113.06
2	A	902	B29	OAE-PBC-OAD	-3.62	104.94	113.06
2	A	902	B29	PBD-CBB-PBC	-3.49	106.56	112.81
2	A	901	B29	OAB-PBD-CBB	3.36	113.69	106.17
2	B	904	B29	OAB-PBD-CBB	3.29	113.55	106.17
2	B	903	B29	CBB-CAS-CAU	-3.16	110.67	116.10
2	A	901	B29	OAA-PBC-CBB	3.00	112.90	106.17
2	A	902	B29	OAA-PBC-CBB	2.83	112.51	106.17
2	A	901	B29	OAE-PBC-OAD	-2.73	106.93	113.06
2	A	902	B29	CAM-CAV-CAW	-2.69	116.34	120.86
2	B	904	B29	OAE-PBC-OAD	-2.67	107.07	113.06
2	A	902	B29	OAG-PBD-OAF	-2.53	107.39	113.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	903	B29	OAG-PBD-OAF	-2.48	107.48	113.06
2	A	902	B29	CAS-CAU-CAL	-2.48	117.75	121.07
2	A	902	B29	OAB-PBD-CBB	2.43	111.63	106.17
2	A	901	B29	OAE-PBC-CBB	-2.32	100.96	106.17
2	B	903	B29	PBD-CBB-PBC	-2.30	108.69	112.81
2	A	902	B29	OAB-PBD-OAG	2.27	114.39	107.99
2	A	901	B29	OAG-PBD-OAF	-2.25	108.00	113.06
2	A	902	B29	OAA-PBC-OAE	2.20	114.20	107.99
2	A	902	B29	CAR-CAV-CAW	2.20	124.30	120.70
2	A	902	B29	CAS-CAU-CAR	2.10	122.90	120.34
2	B	904	B29	OAA-PBC-CBB	2.03	110.72	106.17
2	A	902	B29	PBC-CBB-CAS	2.00	114.62	109.23

There are no chirality outliers.

All (14) torsion outliers are listed below:

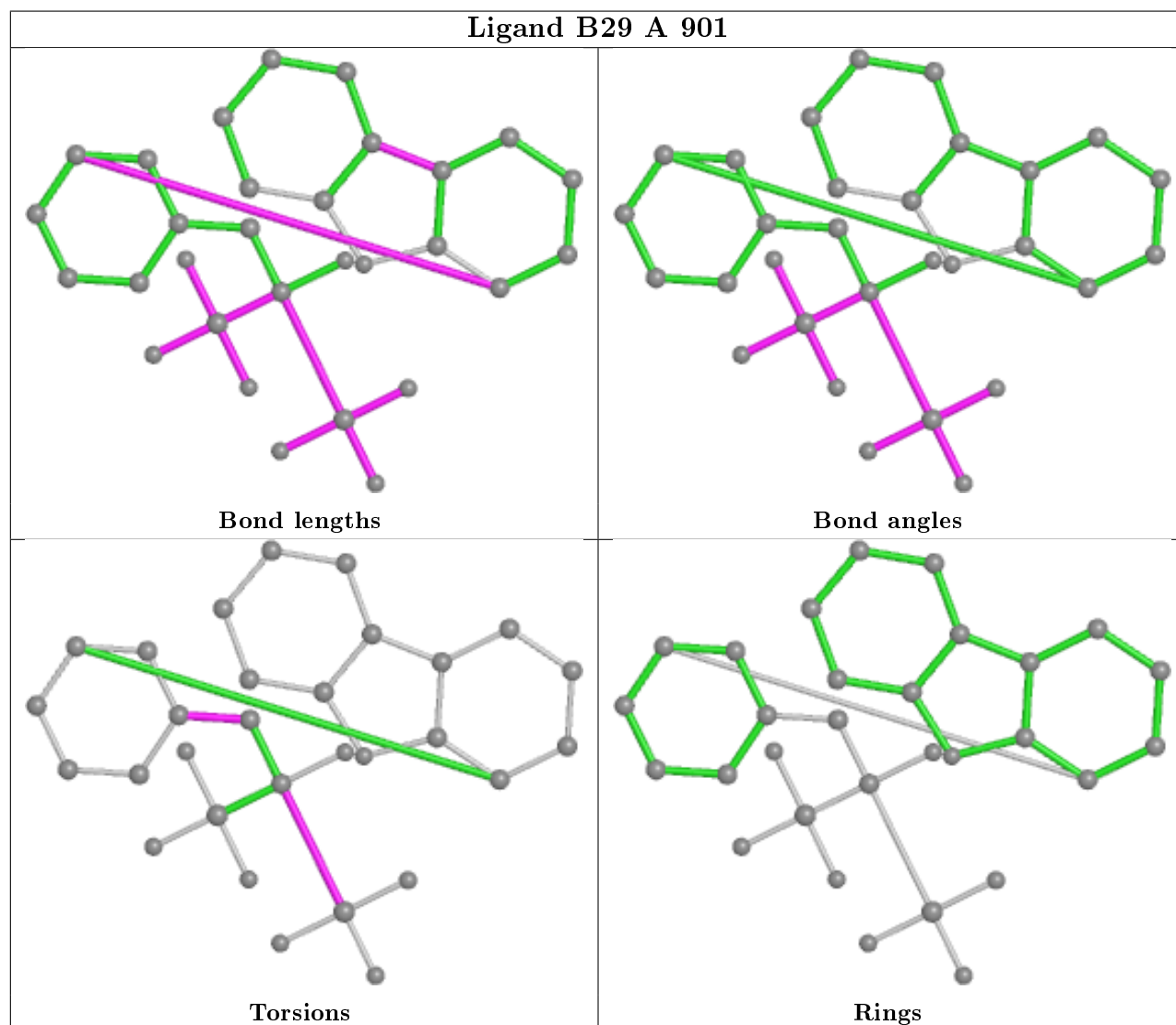
Mol	Chain	Res	Type	Atoms
2	A	902	B29	CAU-CAS-CBB-PBC
2	A	902	B29	CAU-CAS-CBB-PBD
2	A	902	B29	CAU-CAS-CBB-OAC
2	B	903	B29	CAM-CAV-CAW-CAN
2	B	903	B29	CAR-CAV-CAW-CAN
2	B	903	B29	CAM-CAV-CAW-CBA
2	A	901	B29	CBB-CAS-CAU-CAR
2	A	901	B29	CBB-CAS-CAU-CAL
2	A	902	B29	CAM-CAV-CAW-CAN
2	B	903	B29	CAR-CAV-CAW-CBA
2	A	902	B29	CAR-CAV-CAW-CAN
2	B	904	B29	CBB-CAS-CAU-CAR
2	A	901	B29	OAC-CBB-PBD-OAF
2	B	903	B29	OAC-CBB-PBD-OAF

There are no ring outliers.

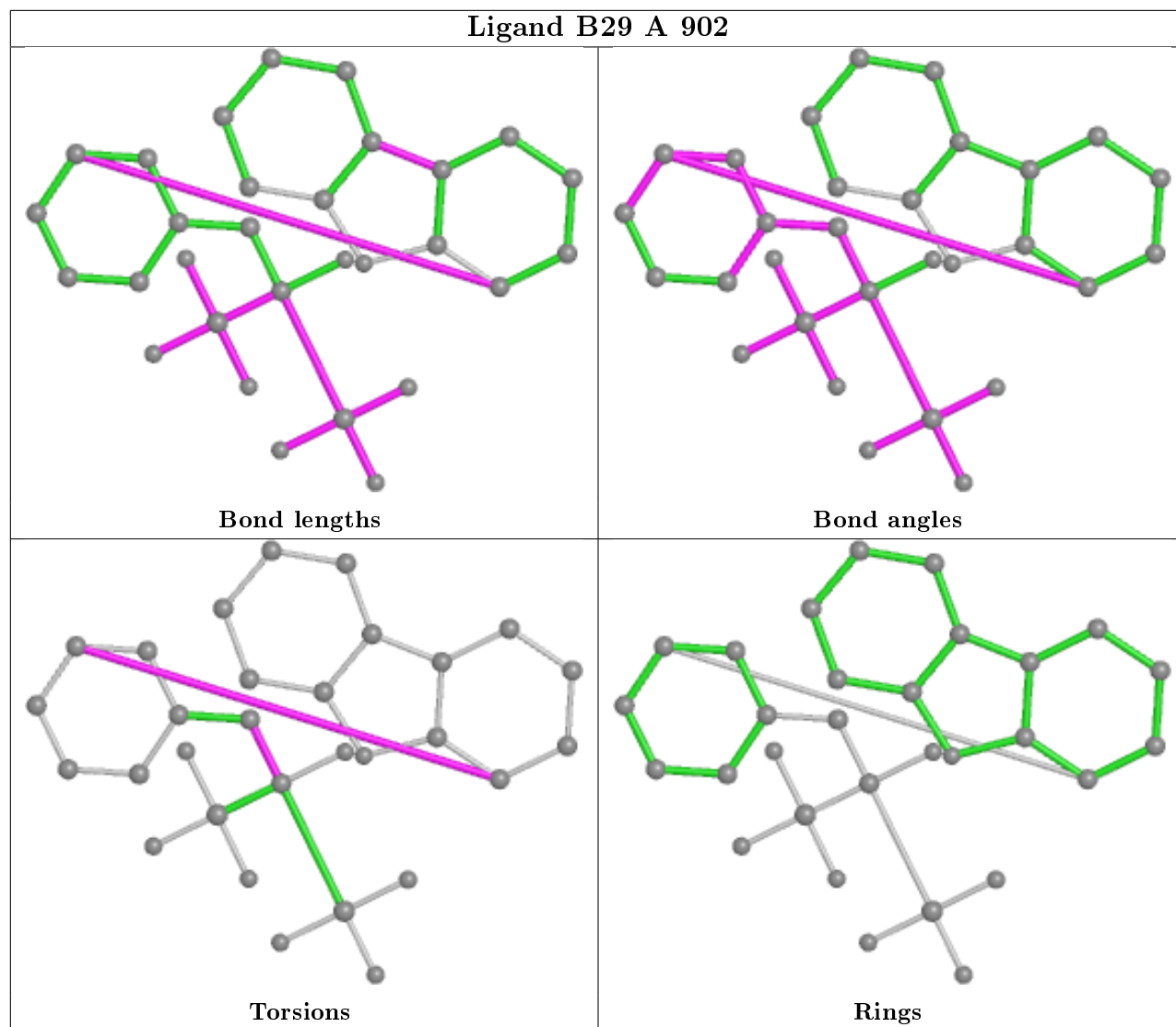
4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	B29	3	0
2	A	902	B29	4	0
2	B	904	B29	6	0
2	B	903	B29	4	0

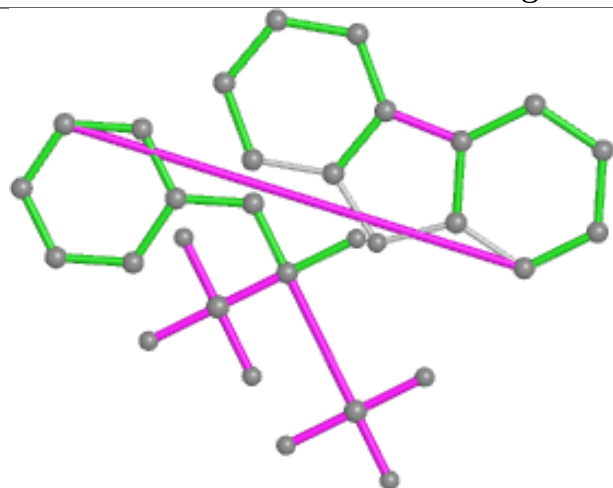
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



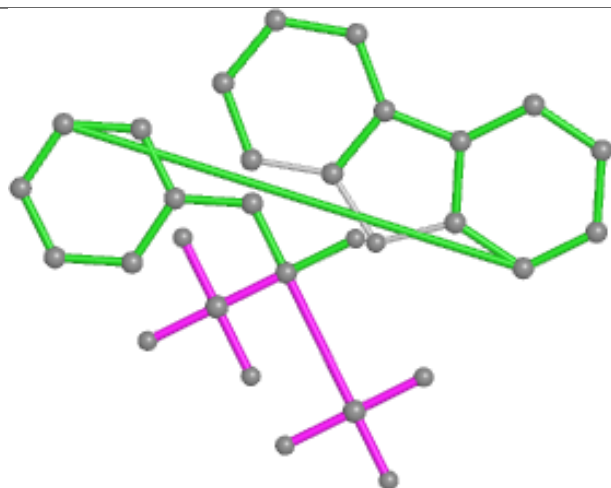
## Ligand B29 A 902



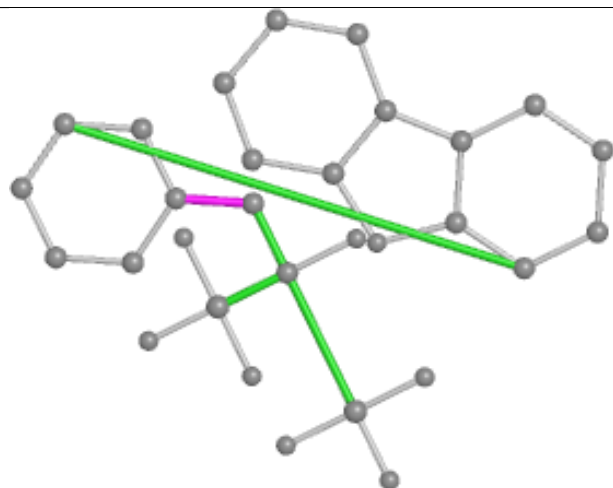
## Ligand B29 B 904



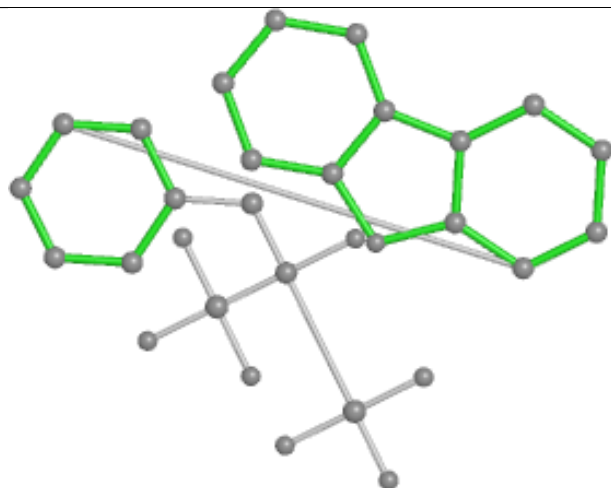
Bond lengths



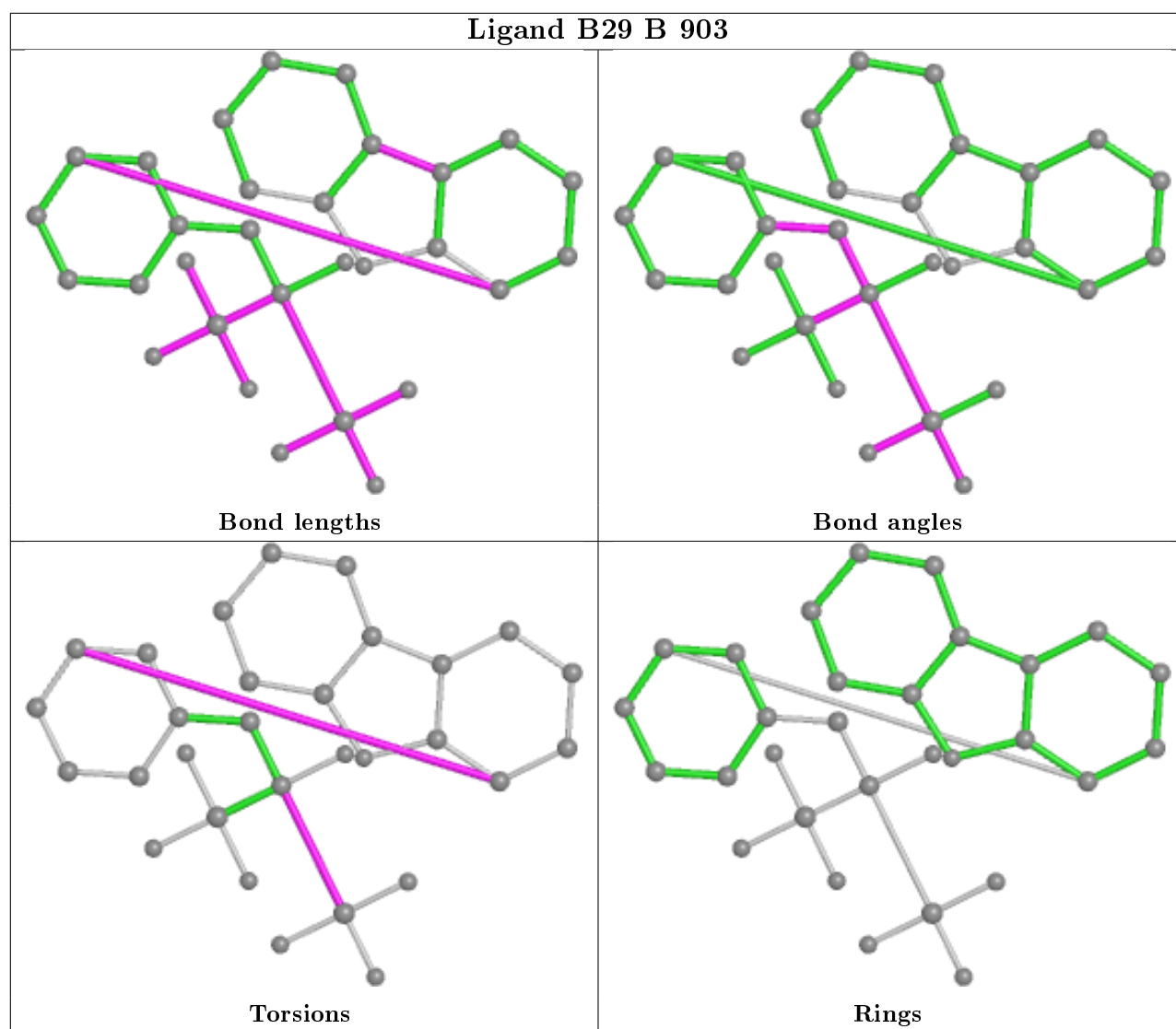
Bond angles



Torsions



Rings



## 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	299/340 (87%)	0.11	17 (5%) 23 28	17, 27, 59, 102	0
1	B	314/340 (92%)	0.28	34 (10%) 5 7	17, 31, 78, 99	0
All	All	613/680 (90%)	0.20	51 (8%) 11 14	17, 29, 67, 102	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	43	PHE	7.5
1	B	332	TYR	6.8
1	B	43	PHE	6.6
1	A	46	ASN	5.6
1	B	197	HIS	5.4
1	A	42	ASN	5.3
1	A	334	ILE	5.1
1	B	230	PHE	4.9
1	B	45	LEU	4.9
1	A	333	ILE	4.2
1	A	332	TYR	4.2
1	B	335	ASP	4.2
1	B	339	GLU	4.2
1	B	309	GLU	3.7
1	B	308	ASN	3.6
1	B	39	PRO	3.4
1	A	197	HIS	3.3
1	A	307	ASP	3.3
1	A	306	ASN	3.3
1	B	329	GLU	3.3
1	B	336	HIS	3.3
1	B	37	LEU	3.2
1	B	4	ASN	3.2
1	B	50	GLN	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	37	LEU	3.1
1	A	336	HIS	3.1
1	A	126	GLU	3.0
1	B	42	ASN	3.0
1	A	39	PRO	2.9
1	B	222	PHE	2.8
1	B	331	LEU	2.8
1	B	338	SER	2.7
1	B	224	MET	2.7
1	B	46	ASN	2.5
1	B	330	LEU	2.4
1	B	126	GLU	2.4
1	B	53	ARG	2.4
1	B	333	ILE	2.4
1	B	38	LYS	2.3
1	A	221	ASP	2.3
1	B	334	ILE	2.3
1	B	41	LYS	2.2
1	A	191	PRO	2.2
1	A	305	LYS	2.2
1	A	124	THR	2.2
1	B	40	GLY	2.2
1	B	310	ASN	2.1
1	B	54	VAL	2.1
1	B	337	LEU	2.1
1	B	191	PRO	2.1
1	B	56	ASN	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,

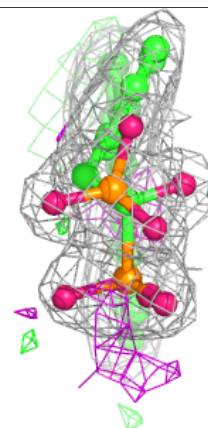
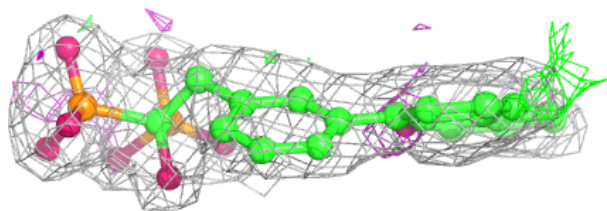
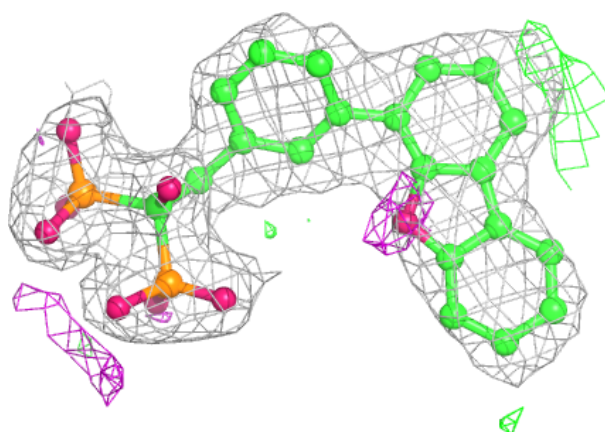
median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	B29	A	902	30/30	0.94	0.13	31,37,44,46	0
2	B29	B	903	30/30	0.95	0.13	34,50,55,56	0
2	B29	B	904	30/30	0.96	0.12	26,30,36,37	0
2	B29	A	901	30/30	0.97	0.12	24,27,30,32	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 B29 A 902:**

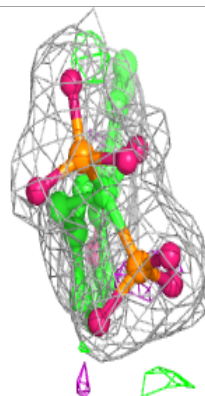
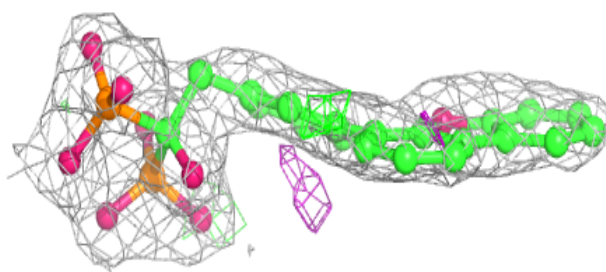
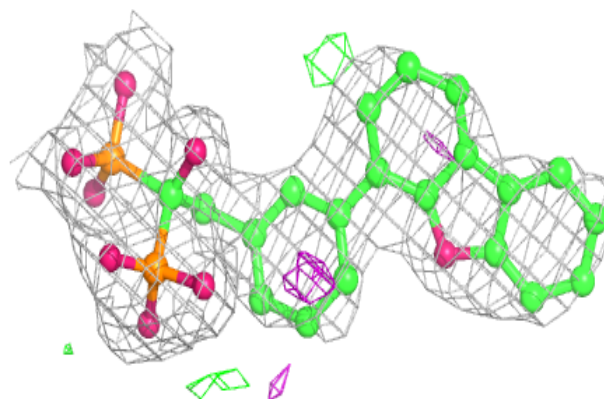
2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



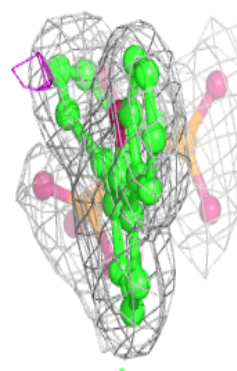
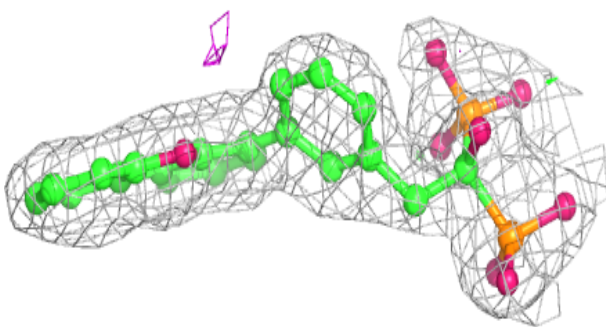
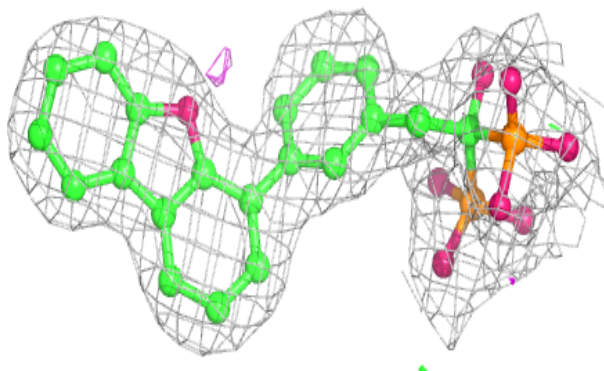


**Electron density around B29 B 903:**

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

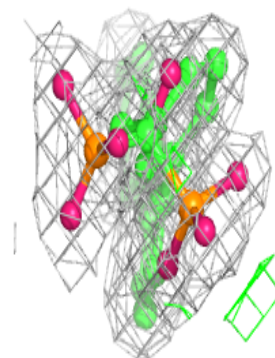
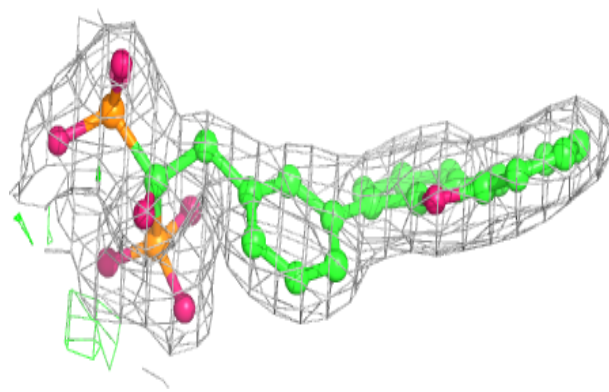
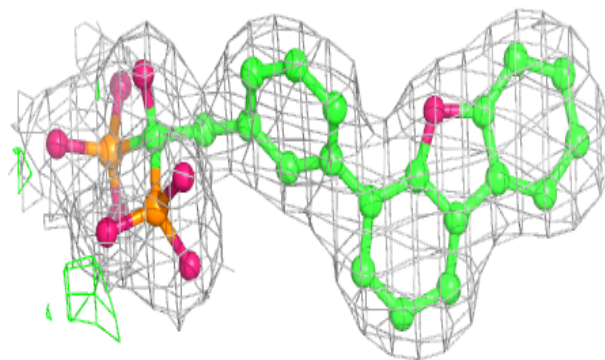
**Electron density around B29 B 904:**

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



**Electron density around B29 A 901:**

$2mF_o-DF_c$  (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.