



# Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2020 – 01:46 am BST

PDB ID : 5D5G  
Title : Structure of colocasia esculenta agglutinin  
Authors : Biswas, H.; Chattopadhyaya, R.  
Deposited on : 2015-08-10  
Resolution : 1.74 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  
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

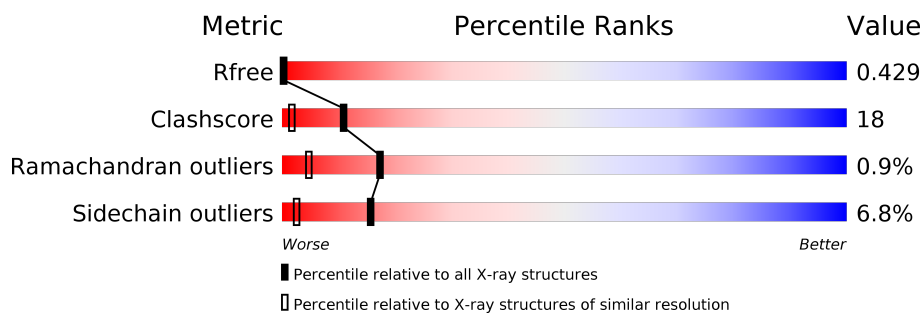
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.74 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	3764 (1.76-1.72)
Clashscore	141614	3923 (1.76-1.72)
Ramachandran outliers	138981	3878 (1.76-1.72)
Sidechain outliers	138945	3878 (1.76-1.72)

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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	109	61% 35% .
1	C	109	59% 39% .
2	B	111	65% 33% ..
2	D	111	59% 38% .

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3577 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 Tuber agglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	109	Total	C	N	O	S	0	0	0
			840	527	147	163	3			
1	C	109	Total	C	N	O	S	0	0	0
			837	524	147	163	3			

- Molecule 2 is a protein called Tuber agglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	110	Total	C	N	O	S	0	0	0
			866	551	149	163	3			
2	D	111	Total	C	N	O	S	0	0	0
			875	556	151	165	3			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Mg	0	0
			3	3		
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	36	Total 36	O 36	0	0
6	B	45	Total 45	O 45	0	0
6	C	22	Total 22	O 22	0	0
6	D	31	Total 31	O 31	0	0



F97	I100	Y101	G102	W106	E107	T108	S109	P110	Q111
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.01 Å 47.20 Å 82.25 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.99 – 1.74 28.85 – 1.54	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.99-1.74) 83.4 (28.85-1.54)	Depositor EDS
$R_{merge}$	0.38	Depositor
$R_{sym}$	0.38	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.25 (at 1.54 Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, $R_{free}$	0.344 , 0.427 0.347 , 0.429	Depositor DCC
$R_{free}$ test set	1021 reflections (1.71%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	6.5	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.19 , -10.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.54$ , $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.79	EDS
Total number of atoms	3577	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.96% 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: MG, EPE, SO4

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	0/858	0.68	0/1165
1	C	0.37	0/855	0.63	0/1161
2	B	0.46	0/889	0.71	0/1203
2	D	0.44	0/898	0.68	0/1215
All	All	0.41	0/3500	0.68	0/4744

There are no bond length outliers.

There are no bond angle outliers.

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	840	0	807	36	0
1	C	837	0	799	34	0
2	B	866	0	824	35	0
2	D	875	0	833	36	1
3	A	2	0	0	0	0
3	B	3	0	0	0	0
4	A	15	0	17	0	0
5	B	5	0	0	0	0
6	A	36	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	45	0	0	7	1
6	C	22	0	0	1	0
6	D	31	0	0	4	2
All	All	3577	0	3280	122	4

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 18.

All (122) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:55:HIS:O	6:B:301:HOH:O	1.98	0.81
1:A:39:ASN:N	1:A:39:ASN:OD1	2.17	0.78
1:A:8:SER:HA	1:A:55:LEU:HB3	1.66	0.77
1:A:11:THR:O	6:A:301:HOH:O	2.05	0.73
1:A:75:ARG:NH1	6:A:302:HOH:O	2.21	0.73
2:B:72:ASP:OD2	6:B:303:HOH:O	2.09	0.69
2:D:52:ASN:OD1	6:D:201:HOH:O	2.09	0.68
2:B:72:ASP:OD1	6:B:304:HOH:O	2.11	0.68
1:A:100:VAL:HG12	2:B:81:SER:HB2	1.77	0.65
2:D:11:SER:HB2	6:D:202:HOH:O	1.95	0.65
1:A:63:ILE:HG13	1:A:72:TRP:HB3	1.80	0.64
1:C:12:LEU:HB3	1:C:53:LEU:HB3	1.81	0.63
1:C:92:ARG:NH1	2:D:107:GLU:OE2	2.32	0.63
1:C:55:LEU:HD13	1:C:61:LEU:HD23	1.82	0.62
2:D:59:ARG:HG2	2:D:60:LEU:H	1.64	0.62
2:D:2:ILE:HB	2:D:13:GLN:HE21	1.65	0.62
2:B:46:GLN:NE2	2:B:48:ASN:HA	2.15	0.61
1:A:106:TRP:NE1	2:B:95:ASP:O	2.31	0.60
2:D:61:ASN:HD21	2:D:65:GLU:HB3	1.64	0.60
1:C:44:THR:HG22	1:C:47:LYS:HD2	1.84	0.60
1:A:88:HIS:CD2	1:A:90:ASP:H	2.21	0.59
1:A:25:LEU:HA	1:A:35:LEU:HA	1.85	0.59
1:A:17:HIS:ND1	1:A:24:ASP:OD1	2.36	0.58
1:A:88:HIS:HD2	1:A:90:ASP:HB2	1.69	0.58
2:B:6:ASN:ND2	6:B:306:HOH:O	2.22	0.57
1:C:8:SER:HA	1:C:55:LEU:HB3	1.86	0.57
1:C:4:ASN:ND2	2:D:7:ASN:OD1	2.37	0.57
2:B:83:SER:OG	2:B:84:LYS:N	2.34	0.57
1:A:14:ARG:NH2	6:A:305:HOH:O	2.37	0.57
2:D:70:ASP:OD1	2:D:74:LYS:N	2.37	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:GLU:HG2	1:C:29:ASP:OD1	2.06	0.56
1:C:28:GLN:OE1	1:C:36:TYR:OH	2.21	0.56
2:B:61:ASN:OD1	2:B:65:GLU:N	2.34	0.56
1:A:44:THR:HA	1:A:47:LYS:HE2	1.88	0.55
2:B:59:ARG:NH2	6:B:307:HOH:O	2.22	0.55
2:B:55:HIS:HB2	2:B:71:ASP:HB2	1.89	0.55
2:B:87:ASP:OD1	2:B:87:ASP:N	2.38	0.55
2:D:28:LEU:HD13	2:D:38:LEU:HB2	1.88	0.55
1:A:32:ASN:ND2	1:A:41:GLN:OE1	2.38	0.55
1:C:83:TYR:CZ	1:C:98:PRO:HG3	2.42	0.55
2:D:85:HIS:NE2	6:D:206:HOH:O	2.33	0.55
2:B:59:ARG:NE	6:B:307:HOH:O	2.35	0.54
1:C:88:HIS:N	1:C:92:ARG:O	2.31	0.53
1:C:44:THR:HG21	1:C:71:VAL:HG12	1.89	0.53
2:D:36:LEU:HB2	2:D:68:ILE:HD11	1.90	0.53
2:B:64:GLY:O	2:B:81:SER:OG	2.22	0.53
2:B:69:LYS:HZ3	2:B:75:THR:HG1	1.56	0.52
2:D:2:ILE:HB	2:D:13:GLN:NE2	2.24	0.52
2:D:69:LYS:NZ	2:D:75:THR:OG1	2.28	0.52
2:D:14:VAL:HG22	2:D:59:ARG:HG3	1.90	0.52
2:B:93:ARG:HB2	2:B:95:ASP:OD1	2.10	0.51
1:C:4:ASN:HD21	2:D:7:ASN:CG	2.13	0.51
1:A:88:HIS:CD2	1:A:90:ASP:HB2	2.47	0.50
2:D:87:ASP:HA	6:D:202:HOH:O	2.11	0.49
1:A:78:SER:OG	1:A:79:VAL:N	2.44	0.49
2:D:93:ARG:HG3	2:D:97:PHE:O	2.13	0.49
2:D:79:SER:O	2:D:80:SER:HB2	2.13	0.49
1:C:17:HIS:HA	1:C:25:LEU:O	2.12	0.49
1:C:2:GLY:O	1:C:3:THR:HG23	2.12	0.49
1:C:3:THR:OG1	1:C:4:ASN:N	2.46	0.49
2:D:8:LEU:O	2:D:21:LEU:HD21	2.12	0.49
2:B:82:SER:O	6:B:305:HOH:O	2.20	0.48
2:B:28:LEU:HD13	2:B:38:LEU:HB2	1.96	0.48
2:B:32:GLY:C	2:B:34:CYS:H	2.17	0.48
2:D:54:GLU:N	2:D:54:GLU:OE2	2.47	0.48
1:C:10:GLN:O	1:C:54:THR:HG22	2.14	0.48
1:A:44:THR:HG21	1:A:72:TRP:HB2	1.96	0.48
1:C:8:SER:O	6:C:201:HOH:O	2.20	0.47
1:C:94:VAL:HG22	2:D:107:GLU:HB3	1.95	0.47
2:D:58:LEU:HD11	2:D:66:LEU:HD21	1.96	0.47
1:C:97:GLY:HA2	2:D:102:GLY:H	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LYS:NZ	1:A:21:GLY:O	2.38	0.47
2:D:49:THR:HB	2:D:52:ASN:ND2	2.29	0.46
2:B:45:TRP:CG	2:B:46:GLN:N	2.83	0.46
1:A:33:LEU:HB3	1:A:42:SER:HB3	1.97	0.46
2:D:36:LEU:HB3	2:D:47:SER:HB3	1.97	0.46
1:C:61:LEU:HB2	1:C:74:SER:HB3	1.97	0.45
2:D:16:TYR:HB3	2:D:57:PHE:HB2	1.98	0.45
1:A:105:PRO:HB2	2:B:41:GLY:HA3	1.98	0.45
1:A:53:LEU:HD13	1:A:63:ILE:HG12	1.97	0.45
2:B:15:LEU:O	2:B:30:MET:HG2	2.16	0.45
2:D:21:LEU:HD23	2:D:21:LEU:HA	1.75	0.45
1:C:23:PHE:HE1	1:C:37:ASN:O	2.00	0.44
1:A:93:LEU:HB3	2:B:108:THR:CG2	2.48	0.44
1:C:76:ALA:HB2	2:D:106:TRP:CD1	2.53	0.44
1:C:78:SER:HB3	1:C:83:TYR:OH	2.18	0.44
1:A:103:ILE:HD13	2:B:45:TRP:CZ3	2.52	0.44
1:C:17:HIS:HD1	1:C:24:ASP:CG	2.21	0.44
2:D:61:ASN:ND2	2:D:65:GLU:HB3	2.32	0.44
1:C:101:PHE:HB3	2:D:100:ILE:HD12	1.99	0.43
1:A:104:ASP:O	1:A:107:VAL:HG12	2.19	0.43
1:A:92:ARG:NE	2:B:109:SER:O	2.43	0.43
1:C:18:LEU:HB2	1:C:25:LEU:HB3	1.99	0.43
1:C:83:TYR:CD2	1:C:97:GLY:HA3	2.53	0.43
1:A:53:LEU:HD12	1:A:53:LEU:HA	1.68	0.43
1:C:18:LEU:HA	1:C:18:LEU:HD23	1.84	0.43
2:B:93:ARG:NH2	2:B:95:ASP:OD2	2.35	0.43
2:B:36:LEU:HB3	2:B:47:SER:HB3	2.00	0.43
1:A:106:TRP:CD1	2:B:97:PHE:HE1	2.37	0.43
2:B:21:LEU:O	2:B:27:GLN:HA	2.19	0.42
1:A:73:ARG:HB2	1:A:75:ARG:HG3	2.01	0.42
1:A:106:TRP:CD1	2:B:97:PHE:CE1	3.08	0.42
1:C:7:LEU:H	1:C:10:GLN:NE2	2.17	0.42
2:D:2:ILE:HD12	2:D:13:GLN:HG3	2.02	0.42
1:A:8:SER:N	1:A:55:LEU:HD23	2.35	0.42
2:B:36:LEU:O	2:B:46:GLN:HA	2.20	0.41
2:D:67:ILE:HG22	2:D:68:ILE:O	2.20	0.41
2:D:36:LEU:HD13	2:D:58:LEU:HD13	2.02	0.41
1:A:4:ASN:N	1:A:4:ASN:OD1	2.52	0.41
2:D:16:TYR:HB3	2:D:57:PHE:CB	2.51	0.41
1:C:5:TYR:OH	2:D:94:ASP:N	2.37	0.41
2:B:69:LYS:HD3	2:B:75:THR:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:LEU:O	2:B:108:THR:HG23	2.21	0.41
1:C:55:LEU:HD13	1:C:61:LEU:CD2	2.49	0.41
1:A:104:ASP:HA	1:A:105:PRO:HD2	1.96	0.41
2:B:66:LEU:HA	2:B:66:LEU:HD12	1.88	0.41
1:C:41:GLN:NE2	1:C:43:ASN:OD1	2.54	0.41
1:A:100:VAL:CG1	2:B:81:SER:HB2	2.49	0.40
2:D:8:LEU:HD22	2:D:91:ILE:HG12	2.01	0.40
1:C:35:LEU:HG	1:C:38:GLY:HA3	2.02	0.40
1:A:1:LEU:N	1:C:15:GLU:HB3	2.36	0.40
1:A:25:LEU:HD11	1:A:33:LEU:HD21	2.03	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:301:HOH:O	6:B:307:HOH:O[4_555]	1.99	0.21
6:D:227:HOH:O	6:D:228:HOH:O[1_545]	2.08	0.12
2:D:62:HIS:ND1	2:D:71:ASP:OD2[4_555]	2.09	0.11
6:D:228:HOH:O	6:D:231:HOH:O[1_565]	2.10	0.10

## 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	107/109 (98%)	93 (87%)	13 (12%)	1 (1%)	17	5
1	C	107/109 (98%)	92 (86%)	13 (12%)	2 (2%)	8	1
2	B	108/111 (97%)	95 (88%)	13 (12%)	0	100	100
2	D	109/111 (98%)	93 (85%)	15 (14%)	1 (1%)	17	5
All	All	431/440 (98%)	373 (86%)	54 (12%)	4 (1%)	17	5

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	3	THR
2	D	50	HIS
1	C	47	LYS
1	A	68	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	91/92 (99%)	82 (90%)	9 (10%)	8	1
1	C	90/92 (98%)	83 (92%)	7 (8%)	12	2
2	B	94/95 (99%)	90 (96%)	4 (4%)	29	8
2	D	95/95 (100%)	90 (95%)	5 (5%)	22	5
All	All	370/374 (99%)	345 (93%)	25 (7%)	16	2

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LEU
1	A	3	THR
1	A	26	VAL
1	A	29	ASP
1	A	39	ASN
1	A	50	ASP
1	A	63	ILE
1	A	69	SER
1	A	73	ARG
2	B	18	ASP
2	B	49	THR
2	B	63	LYS
2	B	87	ASP
1	C	6	LEU
1	C	14	ARG
1	C	44	THR
1	C	73	ARG
1	C	86	VAL

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Mol	Chain	Res	Type
1	C	100	VAL
1	C	107	VAL
2	D	6	ASN
2	D	11	SER
2	D	54	GLU
2	D	70	ASP
2	D	109	SER

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

Mol	Chain	Res	Type
2	B	48	ASN
2	D	52	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 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 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
5	SO4	B	201	-	4,4,4	0.15	0	6,6,6	0.21	0
4	EPE	A	202	-	15,15,15	0.78	1 (6%)	18,20,20	1.76	5 (27%)

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
4	EPE	A	202	-	-	4/9/19/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	202	EPE	C10-S	2.57	1.81	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	202	EPE	C5-N4-C3	4.24	118.36	108.83
4	A	202	EPE	C7-N4-C3	3.12	119.23	111.23
4	A	202	EPE	C7-N4-C5	2.96	118.82	111.23
4	A	202	EPE	O3S-S-C10	2.12	109.20	105.77
4	A	202	EPE	O1S-S-C10	2.12	109.46	106.92

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	202	EPE	C8-C7-N4-C5
4	A	202	EPE	C9-C10-S-O1S
4	A	202	EPE	C9-C10-S-O2S
4	A	202	EPE	C9-C10-S-O3S

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.



## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers ⓘ

Unable to reproduce the depositors R factor - this section is therefore empty.