



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 09:38 AM GMT

PDB ID : 4CC7
Title : Crystal structure of the sixth or C-terminal SH3 domain of human Tuba in complex with proline-rich peptides of N-WASP. Space group P41
Authors : Polle, L.; Rigano, L.; Julian, R.; Ireton, K.; Schubert, W.-D.
Deposited on : 2013-10-18
Resolution : 1.97 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

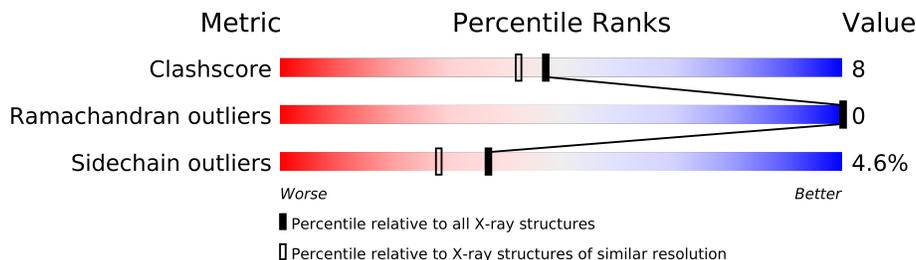
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.97 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))
Clashscore	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	67	
1	C	67	
1	E	67	
1	G	67	
1	I	67	
1	K	67	
1	M	67	
2	B	12	
2	D	12	
2	F	12	
2	H	12	
2	J	12	
2	L	12	
2	N	12	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4418 atoms, of which 0 are hydrogen and 0 are deuterium.

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 DYNAMIN-BINDING PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	59	521	340	91	90	0	10	1
1	C	63	529	343	89	97	0	2	0
1	E	62	546	354	94	98	0	7	0
1	G	61	505	330	84	91	0	1	0
1	I	61	515	333	89	93	0	2	0
1	K	61	534	350	91	93	0	10	1
1	M	60	502	330	84	88	0	4	1

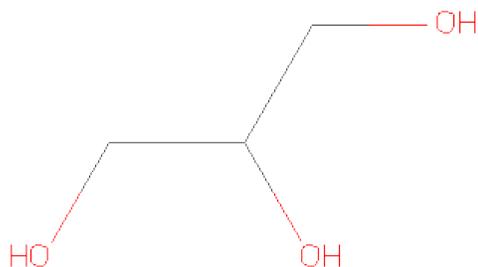
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
A	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
C	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
C	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
E	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
E	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
G	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
G	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
I	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
I	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
K	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
K	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7
M	1511	GLY	-	EXPRESSION TAG	UNP Q6XZF7
M	1512	PRO	-	EXPRESSION TAG	UNP Q6XZF7

- Molecule 2 is a protein called NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	11	Total 68	C 44	N 11	O 13	0	1	1
2	D	12	Total 73	C 47	N 12	O 14	0	0	1
2	F	12	Total 73	C 47	N 12	O 14	0	0	1
2	H	12	Total 75	C 48	N 12	O 15	0	0	0
2	J	12	Total 72	C 46	N 12	O 14	0	0	1
2	L	12	Total 68	C 43	N 11	O 14	0	0	1
2	N	12	Total 73	C 47	N 12	O 14	0	0	1

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	Total 6	C 3	O 3	0	0
3	C	1	Total 6	C 3	O 3	0	0
3	K	1	Total 6	C 3	O 3	0	0
3	M	1	Total 6	C 3	O 3	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	G	2	Total Cl 2 2	0	0
4	K	3	Total Cl 3 3	0	0
4	E	2	Total Cl 2 2	0	0
4	I	1	Total Cl 1 1	0	0
4	C	3	Total Cl 3 3	0	0
4	A	2	Total Cl 2 2	0	0
4	M	4	Total Cl 4 4	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	K	1	Total Mg 1 1	0	0
5	M	1	Total Mg 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	20	Total O 20 20	0	0
6	C	44	Total O 44 44	0	0
6	D	1	Total O 1 1	0	0
6	E	29	Total O 29 29	0	0
6	F	4	Total O 4 4	0	0
6	G	30	Total O 30 30	0	0
6	H	1	Total O 1 1	0	0
6	I	21	Total O 21 21	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	K	31	Total 31	O 31	0	0
6	L	4	Total 4	O 4	0	0
6	M	31	Total 31	O 31	0	0
6	N	5	Total 5	O 5	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain A: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain C: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain E: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain G: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain I: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain K: 



- Molecule 1: DYNAMIN-BINDING PROTEIN

Chain M:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain B:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain D:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain F:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain H:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain J:

There are no outlier residues recorded for this chain.

- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain L:



- Molecule 2: NEURAL WISKOTT-ALDRICH SYNDROME PROTEIN

Chain N:

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	88.57Å 88.57Å 69.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.59 – 1.97	Depositor
% Data completeness (in resolution range)	99.8 (46.59-1.97)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.176 , 0.215	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	4418	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

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, MG, CL

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	0.94	0/579	1.06	1/780 (0.1%)
1	C	1.10	1/553 (0.2%)	1.03	0/747
1	E	0.88	0/585	0.92	0/789
1	G	0.91	0/523	0.89	0/707
1	I	0.95	1/535 (0.2%)	1.08	7/722 (1.0%)
1	K	1.18	0/589	1.06	2/798 (0.3%)
1	M	0.96	0/531	0.98	1/719 (0.1%)
2	B	0.80	0/76	0.73	0/107
2	D	1.44	1/77 (1.3%)	0.82	0/108
2	F	1.14	1/77 (1.3%)	1.02	0/108
2	H	0.83	0/79	0.69	0/110
2	J	0.70	0/76	0.71	0/107
2	L	0.96	0/71	0.71	0/100
2	N	1.00	0/77	0.79	0/108
All	All	1.00	4/4428 (0.1%)	0.98	11/6010 (0.2%)

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

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	12	GLY	N-CA	9.86	1.60	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	1518	TYR	CE2-CZ	-8.23	1.27	1.38
1	C	1544	GLU	CD-OE2	-6.68	1.18	1.25
2	F	12	GLY	N-CA	5.84	1.54	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1540	LEU	CB-CG-CD1	-13.13	88.67	111.00
1	I	1518	TYR	OH-CZ-CE2	-6.97	101.27	120.10
1	I	1572[A]	ARG	NE-CZ-NH2	-6.83	116.88	120.30
1	I	1572[B]	ARG	NE-CZ-NH2	-6.83	116.88	120.30
1	I	1518	TYR	CE1-CZ-OH	6.58	137.87	120.10
1	I	1572[A]	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	I	1572[B]	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	K	1572	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	K	1527	ARG	CG-CD-NE	5.38	123.10	111.80
1	I	1530	ASN	CB-CA-C	-5.24	99.93	110.40
1	M	1546	LYS	CB-CG-CD	5.02	124.66	111.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	11	SER	Peptide
2	F	11	SER	Peptide
1	K	1515	ASN	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	521	0	0	6	0
1	C	529	0	0	4	0
1	E	546	0	0	4	1
1	G	505	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	I	515	0	0	3	0
1	K	534	0	0	10	0
1	M	502	0	0	2	0
2	B	68	0	0	0	0
2	D	73	0	0	2	0
2	F	73	0	0	0	0
2	H	75	0	0	0	0
2	J	72	0	0	0	0
2	L	68	0	0	1	0
2	N	73	0	0	0	0
3	A	6	0	8	0	0
3	C	6	0	8	3	0
3	K	6	0	8	1	0
3	M	6	0	8	0	0
4	A	2	0	0	0	0
4	C	3	0	0	0	0
4	E	2	0	0	1	0
4	G	2	0	0	0	0
4	I	1	0	0	0	0
4	K	3	0	0	0	0
4	M	4	0	0	0	0
5	K	1	0	0	0	0
5	M	1	0	0	0	0
6	A	20	0	0	3	0
6	C	44	0	0	3	0
6	D	1	0	0	1	0
6	E	29	0	0	0	0
6	F	4	0	0	0	0
6	G	30	0	0	1	0
6	H	1	0	0	0	0
6	I	21	0	0	2	1
6	K	31	0	0	4	0
6	L	4	0	0	0	0
6	M	31	0	0	1	0
6	N	5	0	0	0	0
All	All	4418	0	32	34	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (34) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1527[A]:ARG:NH2	6:A:2002:HOH:O	1.78	1.15
2:D:12:GLY:CA	2:L:12:GLY:N	2.50	0.74
1:A:1519:PHE:CE2	1:A:1521:VAL:CG1	2.74	0.71
3:C:2578:GOL:C3	6:C:2044:HOH:O	2.47	0.62
1:E:1515[A]:ASN:CG	1:E:1516[A]:GLN:N	2.53	0.61
1:M:1575:GLU:N	6:M:2031:HOH:O	2.36	0.59
1:A:1560:ASN:OD1	6:A:2016:HOH:O	2.16	0.58
1:E:1515[A]:ASN:OD1	1:E:1573:LYS:NZ	2.40	0.54
1:A:1553:GLU:OE1	6:A:2013:HOH:O	2.19	0.54
1:C:1515:ASN:OD1	1:C:1573:LYS:NZ	2.41	0.52
1:E:1515[B]:ASN:O	1:E:1516[B]:GLN:CB	2.57	0.52
1:G:1538:GLN:NE2	6:G:2015:HOH:O	2.43	0.52
6:C:2019:HOH:O	1:M:1541:LYS:NZ	2.42	0.51
6:D:2001:HOH:O	1:K:1528[A]:ASN:ND2	2.44	0.50
1:E:1562[B]:LYS:NZ	4:E:2578:CL:CL	2.82	0.50
1:K:1569[B]:ASN:ND2	6:K:2024:HOH:O	2.45	0.50
1:C:1519:PHE:CD2	1:C:1575:GLU:OE2	2.66	0.48
1:K:1566[B]:VAL:CG2	1:K:1571[B]:ILE:CD1	2.92	0.47
1:I:1571:ILE:O	1:I:1572[A]:ARG:NE	2.47	0.47
1:K:1562[A]:LYS:NZ	6:K:2011:HOH:O	2.47	0.47
1:I:1515:ASN:N	6:I:2002:HOH:O	2.47	0.47
3:C:2578:GOL:H32	6:C:2044:HOH:O	2.11	0.45
1:K:1528[A]:ASN:ND2	6:K:2009:HOH:O	2.48	0.45
1:C:1573:LYS:N	3:C:2578:GOL:HO3	2.15	0.45
1:K:1516[A]:GLN:CG	1:K:1573[A]:LYS:NZ	2.81	0.44
1:I:1515:ASN:CB	6:I:2002:HOH:O	2.67	0.43
1:C:1519:PHE:CD2	1:C:1575:GLU:CD	2.92	0.42
1:K:1537:ASN:ND2	6:K:2003:HOH:O	2.52	0.42
2:D:1:PRO:N	2:D:2:PRO:CD	2.83	0.41
1:K:1528[A]:ASN:ND2	1:K:1530:ASN:ND2	2.67	0.41
1:K:1535:SER:OG	3:K:2576:GOL:H12	2.20	0.40

All (1) 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	Distance(Å)	Clash(Å)
1:E:1546:LYS:NZ	6:I:2002:HOH:O[4.644]	2.14	0.06

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	67/67 (100%)	67 (100%)	0	0	100	100
1	C	63/67 (94%)	63 (100%)	0	0	100	100
1	E	66/67 (98%)	64 (97%)	2 (3%)	0	100	100
1	G	60/67 (90%)	60 (100%)	0	0	100	100
1	I	61/67 (91%)	61 (100%)	0	0	100	100
1	K	69/67 (103%)	67 (97%)	2 (3%)	0	100	100
1	M	62/67 (92%)	62 (100%)	0	0	100	100
2	B	10/12 (83%)	10 (100%)	0	0	100	100
2	D	10/12 (83%)	10 (100%)	0	0	100	100
2	F	10/12 (83%)	10 (100%)	0	0	100	100
2	H	10/12 (83%)	10 (100%)	0	0	100	100
2	J	10/12 (83%)	10 (100%)	0	0	100	100
2	L	10/12 (83%)	10 (100%)	0	0	100	100
2	N	10/12 (83%)	10 (100%)	0	0	100	100
All	All	518/553 (94%)	514 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	61/58 (105%)	57 (93%)	4 (7%)	24	15
1	C	58/58 (100%)	55 (95%)	3 (5%)	32	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	62/58 (107%)	57 (92%)	5 (8%)	17	9
1	G	54/58 (93%)	51 (94%)	3 (6%)	30	20
1	I	56/58 (97%)	55 (98%)	1 (2%)	71	71
1	K	63/58 (109%)	59 (94%)	4 (6%)	25	16
1	M	56/58 (97%)	51 (91%)	5 (9%)	14	7
2	B	9/9 (100%)	9 (100%)	0	100	100
2	D	9/9 (100%)	9 (100%)	0	100	100
2	F	9/9 (100%)	9 (100%)	0	100	100
2	H	9/9 (100%)	8 (89%)	1 (11%)	9	4
2	J	9/9 (100%)	9 (100%)	0	100	100
2	L	8/9 (89%)	8 (100%)	0	100	100
2	N	9/9 (100%)	9 (100%)	0	100	100
All	All	472/469 (101%)	446 (94%)	26 (6%)	37	21

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1530	ASN
1	A	1540	LEU
1	A	1569[A]	ASN
1	A	1569[B]	ASN
1	C	1530	ASN
1	C	1538	GLN
1	C	1569	ASN
1	E	1515[A]	ASN
1	E	1515[B]	ASN
1	E	1528	ASN
1	E	1562[A]	LYS
1	E	1562[B]	LYS
1	G	1530	ASN
1	G	1562[A]	LYS
1	G	1562[B]	LYS
2	H	8	SER
1	I	1546	LYS
1	K	1516[A]	GLN
1	K	1516[B]	GLN
1	K	1530	ASN
1	K	1574	THR

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Mol	Chain	Res	Type
1	M	1540[A]	LEU
1	M	1540[B]	LEU
1	M	1546	LYS
1	M	1563	LYS
1	M	1569	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 23 ligands modelled in this entry, 19 are monoatomic - leaving 4 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	$\# Z > 2$
3	GOL	A	2575	-	5,5,5	0.52	0	5,5,5	2.02	2 (40%)
3	GOL	C	2578	-	5,5,5	0.53	0	5,5,5	1.15	1 (20%)
3	GOL	K	2576	-	5,5,5	0.51	0	5,5,5	0.55	0
3	GOL	M	2576	-	5,5,5	0.42	0	5,5,5	0.39	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
3	GOL	A	2575	-	-	0/4/4/4	0/0/0/0
3	GOL	C	2578	-	-	0/4/4/4	0/0/0/0
3	GOL	K	2576	-	-	0/4/4/4	0/0/0/0
3	GOL	M	2576	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2575	GOL	O3-C3-C2	-3.58	92.23	109.71
3	A	2575	GOL	O2-C2-C1	2.36	118.97	108.22
3	C	2578	GOL	O3-C3-C2	2.19	120.40	109.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.