



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 05:56 am BST

PDB ID : 3LS3
Title : Padron0.9-ON (fluorescent state)
Authors : Brakemann, T.; Weber, G.; Trowitzsch, S.; Wahl, M.C.; Jakobs, S.
Deposited on : 2010-02-12
Resolution : 1.65 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

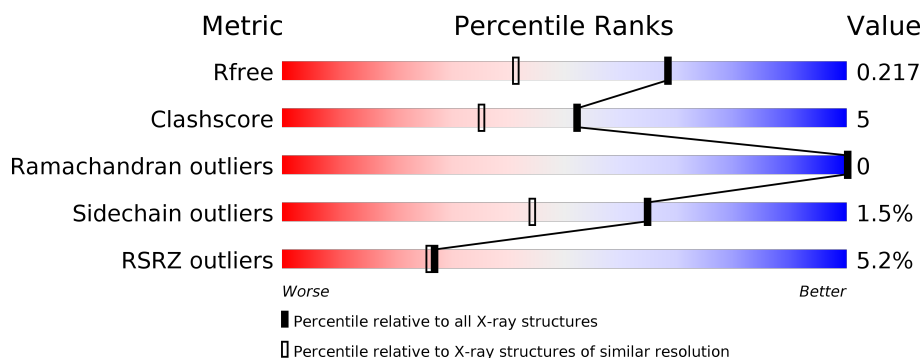
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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 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	235	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>9%</div> <div>8%</div> </div> </div>
1	B	235	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>9%</div> </div> </div>
1	C	235	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>9%</div> </div> </div>
1	D	235	<div> <div>5%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div></div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SPD	B	225	-	-	X	-

2 Entry composition [i](#)

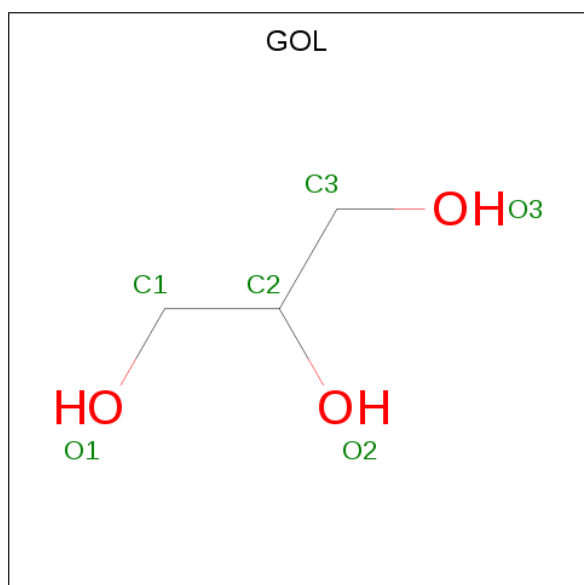
There are 5 unique types of molecules in this entry. The entry contains 8231 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Padron0.9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	16	0
			1846	1172	311	348	15			
1	B	215	Total	C	N	O	S	0	19	0
			1857	1188	312	344	13			
1	C	214	Total	C	N	O	S	0	16	0
			1816	1159	309	335	13			
1	D	225	Total	C	N	O	S	0	18	0
			1921	1223	329	355	14			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



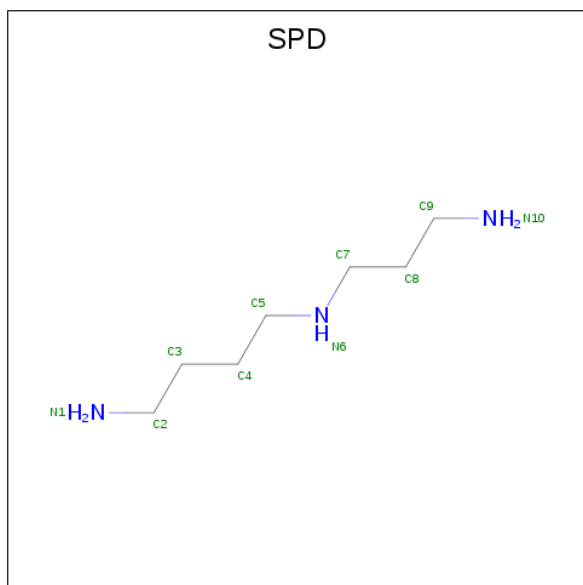
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SPERMIDINE (three-letter code: SPD) (formula: $C_7H_{19}N_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			10	7	3		
3	B	1	Total	C	N	0	0
			10	7	3		
3	B	1	Total	C	N	0	0
			10	7	3		
3	C	1	Total	C	N	0	0
			10	7	3		
3	C	1	Total	C	N	0	0
			10	7	3		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).

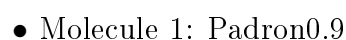
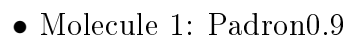
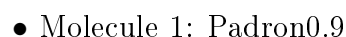


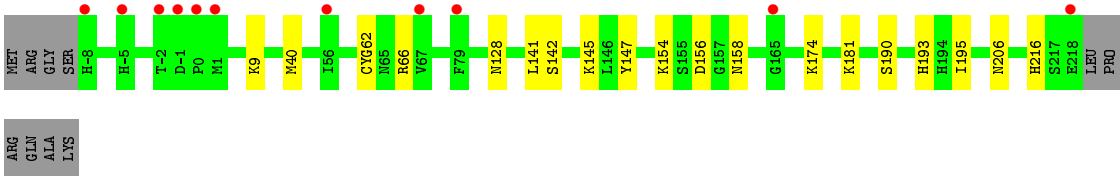
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	156	Total	O	0	10
			166	166		
5	B	169	Total	O	0	1
			170	170		
5	C	153	Total	O	0	2
			155	155		
5	D	202	Total	O	0	3
			205	205		

- Molecule 1: Padron0.9





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.86Å 103.81Å 121.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.65 19.97 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-1.65) 99.3 (19.97-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.66Å)	Xtriage
Refinement program	REFMAC 5.5.0063	Depositor
R, R_{free}	0.168 , 0.200 0.183 , 0.217	Depositor DCC
R_{free} test set	5479 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8231	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, EPE, GYC, SPD

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	0.55	0/1888	0.64	0/2544
1	B	0.66	0/1911	0.69	0/2570
1	C	0.65	0/1877	0.73	0/2528
1	D	0.71	0/1992	0.74	0/2685
All	All	0.65	0/7668	0.70	0/10327

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	1846	0	1771	29	0
1	B	1857	0	1823	17	0
1	C	1816	0	1771	16	0
1	D	1921	0	1854	24	0
2	A	12	0	16	0	0
2	C	6	0	8	0	0
2	D	12	0	16	1	0
3	A	10	0	19	1	0
3	B	20	0	38	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	20	0	38	2	0
4	D	15	0	17	0	0
5	A	166	0	0	4	0
5	B	170	0	0	2	0
5	C	155	0	0	2	0
5	D	205	0	0	9	0
All	All	8231	0	7371	82	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 5.

All (82) 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:D:158[A]:ASN:OD1	5:D:508:HOH:O	1.55	1.24
1:C:158[A]:ASN:OD1	5:C:637:HOH:O	1.54	1.19
1:B:59[B]:MET:HG2	5:B:250:HOH:O	1.39	1.19
1:A:135[B]:ARG:HG2	1:A:135[B]:ARG:HH11	1.08	1.13
1:A:135[B]:ARG:CG	1:A:135[B]:ARG:HH11	1.76	0.98
1:D:206:ASN:HB3	2:D:227:GOL:H11	1.54	0.89
1:A:135[B]:ARG:NH1	5:A:654[B]:HOH:O	2.07	0.88
1:A:135[B]:ARG:HG2	1:A:135[B]:ARG:NH1	1.90	0.86
1:A:73:GLU:H	1:A:73:GLU:CD	1.87	0.77
1:C:170[B]:ARG:HD2	1:D:147:TYR:CZ	2.23	0.72
1:A:135[B]:ARG:CG	1:A:135[B]:ARG:NH1	2.45	0.72
1:C:156[A]:ASP:OD1	1:C:174:LYS:HD3	1.92	0.69
1:D:216[A]:HIS:HD2	5:D:488:HOH:O	1.75	0.69
1:A:12:LEU:HD23	1:A:40[B]:MET:CE	2.23	0.68
1:A:135[B]:ARG:HG2	5:A:654[B]:HOH:O	1.94	0.67
1:A:174:LYS:HZ3	3:B:225:SPD:HN6	1.44	0.66
1:D:190[A]:SER:HG	1:D:216[A]:HIS:CD2	2.13	0.65
1:A:12:LEU:HD23	1:A:40[B]:MET:HE1	1.77	0.65
1:A:10:ILE:HG21	1:A:40[B]:MET:CE	2.28	0.64
1:B:107[A]:ILE:HD12	1:B:107[A]:ILE:N	2.14	0.61
1:A:10:ILE:HG21	1:A:40[B]:MET:HE2	1.82	0.61
1:D:156[A]:ASP:OD1	1:D:174:LYS:CD	2.49	0.61
1:C:5:LYS:H	1:C:8[B]:MET:HE2	1.67	0.60
1:A:62:GYC:HE2	1:A:195:ILE:HB	1.84	0.59
1:B:59[B]:MET:CG	5:B:250:HOH:O	2.18	0.58
1:D:156[A]:ASP:OD1	1:D:174:LYS:HD2	2.03	0.57
1:B:156[B]:ASP:OD1	1:B:174[B]:LYS:HG2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:LYS:HE3	1:B:112:ASP:HB3	1.87	0.57
1:C:174:LYS:CE	5:D:599:HOH:O	2.53	0.56
1:B:190[B]:SER:OG	1:B:216:HIS:CE1	2.60	0.54
1:D:156[B]:ASP:HB3	1:D:174:LYS:HD3	1.89	0.54
1:C:134:LYS:HE3	3:C:226:SPD:H72	1.91	0.52
1:A:135[B]:ARG:CB	1:A:135[B]:ARG:NH1	2.72	0.52
1:D:190[A]:SER:OG	1:D:216[A]:HIS:CD2	2.63	0.52
1:D:190[A]:SER:HG	1:D:216[A]:HIS:CE1	2.27	0.51
1:A:5:LYS:HG3	1:A:112[B]:ASP:OD2	2.10	0.51
1:A:174:LYS:NZ	3:B:225:SPD:HN6	2.08	0.51
1:B:62:GYC:HE2	1:B:195:ILE:HB	1.93	0.49
1:A:142[B]:SER:HB2	1:A:159:TYR:CE1	2.48	0.49
1:D:128:ASN:HB3	5:D:444:HOH:O	2.13	0.49
1:C:62:GYC:HE1	1:C:195:ILE:HB	1.96	0.48
1:B:170:ARG:CB	3:B:225:SPD:H71	2.44	0.47
1:A:31:GLY:HA3	1:A:68:PHE:CE1	2.50	0.47
1:B:96:GLU:HB3	3:B:225:SPD:H72	1.95	0.47
1:C:43[B]:LYS:HG2	1:C:206:ASN:OD1	2.14	0.47
1:A:12:LEU:HD23	1:A:40[B]:MET:HE3	1.96	0.47
1:C:170[B]:ARG:HD2	1:D:147:TYR:CE1	2.50	0.47
1:D:40[B]:MET:HE3	1:D:40[B]:MET:HB3	1.51	0.47
1:A:10:ILE:HG21	1:A:40[B]:MET:HE1	1.96	0.47
1:D:62:GYC:HE2	1:D:195:ILE:HB	1.97	0.46
1:A:73:GLU:N	1:A:73:GLU:CD	2.62	0.46
1:D:40[B]:MET:HE2	1:D:40[B]:MET:HB2	1.39	0.46
1:A:1:MET:HB3	1:A:4:ILE:H	1.81	0.45
1:A:123:VAL:HB	1:C:90:GLU:HB3	1.99	0.45
1:C:174:LYS:HE2	5:D:599:HOH:O	2.15	0.45
1:D:9[B]:LYS:HB2	1:D:9[B]:LYS:HE3	1.73	0.45
1:C:97:ASP:HB3	3:C:227:SPD:H31	1.98	0.44
5:C:246:HOH:O	1:D:145:LYS:HE3	2.17	0.44
1:C:158[B]:ASN:ND2	1:D:158[B]:ASN:OD1	2.51	0.44
1:A:94:ILE:HG22	3:A:227:SPD:H92	1.99	0.43
1:B:170:ARG:HB2	3:B:225:SPD:H71	1.99	0.43
1:B:59[B]:MET:HG3	1:B:173:PHE:CZ	2.54	0.43
1:D:156[A]:ASP:OD1	1:D:174:LYS:HD3	2.19	0.43
1:C:43[A]:LYS:HE2	1:C:43[A]:LYS:HB2	1.67	0.43
1:B:142[B]:SER:HB2	1:B:159:TYR:CE1	2.55	0.42
1:A:10:ILE:CG2	1:A:40[B]:MET:HE1	2.49	0.42
1:C:148:VAL:HG21	1:C:185:LEU:HB3	2.01	0.42
1:C:145:LYS:HE3	5:D:236:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:158[A]:ASN:ND2	5:D:647:HOH:O	2.52	0.42
1:D:142[A]:SER:OG	1:D:193:HIS:HB2	2.20	0.42
1:A:62:GYC:N2	1:A:62:GYC:HD2	2.35	0.42
1:A:142[A]:SER:HB3	1:A:159:TYR:CE1	2.54	0.41
1:D:190[A]:SER:OG	1:D:216[A]:HIS:NE2	2.44	0.41
5:A:452:HOH:O	1:B:145:LYS:HE3	2.19	0.41
1:D:154:LYS:NZ	5:D:375:HOH:O	2.54	0.41
5:A:635:HOH:O	3:B:225:SPD:H51	2.20	0.41
1:B:59[B]:MET:H	1:B:59[B]:MET:HG2	1.65	0.41
1:D:174:LYS:HE2	5:D:668:HOH:O	2.20	0.41
1:B:62:GYC:HD2	1:B:62:GYC:N2	2.35	0.40
1:B:66:ARG:HB2	1:B:66:ARG:CZ	2.51	0.40
1:A:45:LYS:HD2	1:A:45:LYS:HA	1.91	0.40
1:A:156[B]:ASP:OD1	1:B:170:ARG:NE	2.48	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	229/235 (97%)	228 (100%)	1 (0%)	0	100	100
1	B	232/235 (99%)	230 (99%)	2 (1%)	0	100	100
1	C	227/235 (97%)	226 (100%)	1 (0%)	0	100	100
1	D	240/235 (102%)	239 (100%)	1 (0%)	0	100	100
All	All	928/940 (99%)	923 (100%)	5 (0%)	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	200/202 (99%)	196 (98%)	4 (2%)	55	32
1	B	204/202 (101%)	202 (99%)	2 (1%)	76	62
1	C	199/202 (98%)	197 (99%)	2 (1%)	76	62
1	D	212/202 (105%)	209 (99%)	3 (1%)	67	46
All	All	815/808 (101%)	804 (99%)	11 (1%)	65	50

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	67	VAL
1	A	73	GLU
1	A	211	GLU
1	B	193	HIS
1	B	218	GLU
1	C	66	ARG
1	C	193	HIS
1	D	66	ARG
1	D	141	LEU
1	D	181	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	216	HIS
1	D	-8	HIS
1	D	21	HIS
1	D	38	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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$
1	GYC	C	62	1	22,22,23	2.75	6 (27%)	26,30,32	2.42	7 (26%)
1	GYC	B	62	1	22,22,23	3.06	6 (27%)	26,30,32	2.91	9 (34%)
1	GYC	D	62	1	22,22,23	2.95	5 (22%)	26,30,32	2.49	8 (30%)
1	GYC	A	62	1	22,22,23	2.73	5 (22%)	26,30,32	2.67	7 (26%)

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
1	GYC	C	62	1	-	2/9/29/30	0/2/2/2
1	GYC	B	62	1	-	2/9/29/30	0/2/2/2
1	GYC	D	62	1	-	1/9/29/30	0/2/2/2
1	GYC	A	62	1	-	2/9/29/30	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	62	GYC	CA2-C2	-9.09	1.39	1.48
1	D	62	GYC	CB2-CA2	8.94	1.42	1.35
1	A	62	GYC	CA2-C2	-7.97	1.40	1.48
1	C	62	GYC	CB2-CA2	7.61	1.41	1.35
1	B	62	GYC	CB2-CA2	7.51	1.41	1.35
1	C	62	GYC	CA2-C2	-6.72	1.42	1.48
1	D	62	GYC	CA2-C2	-6.63	1.42	1.48
1	A	62	GYC	CB2-CA2	6.58	1.40	1.35
1	B	62	GYC	OH-CZ	-5.75	1.23	1.37
1	A	62	GYC	OH-CZ	-5.37	1.24	1.37
1	D	62	GYC	OH-CZ	-5.36	1.24	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	62	GYC	OH-CZ	-5.18	1.24	1.37
1	A	62	GYC	CG2-CB2	-3.18	1.40	1.46
1	C	62	GYC	C2-N3	-3.11	1.32	1.39
1	B	62	GYC	C2-N3	-3.01	1.32	1.39
1	D	62	GYC	C2-N3	-3.00	1.32	1.39
1	D	62	GYC	CG2-CB2	-2.82	1.41	1.46
1	A	62	GYC	C2-N3	-2.72	1.33	1.39
1	B	62	GYC	CG2-CB2	-2.52	1.42	1.46
1	B	62	GYC	C1-N3	-2.39	1.33	1.37
1	C	62	GYC	CG2-CB2	-2.31	1.42	1.46
1	C	62	GYC	CB1-CA1	2.11	1.55	1.53

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	GYC	CA2-C2-N3	9.73	107.97	103.37
1	B	62	GYC	CA2-C2-N3	9.71	107.96	103.37
1	C	62	GYC	CA2-C2-N3	9.30	107.77	103.37
1	D	62	GYC	CA2-C2-N3	8.45	107.37	103.37
1	B	62	GYC	O2-C2-CA2	-5.83	127.68	130.96
1	A	62	GYC	CG2-CB2-CA2	-4.91	123.93	129.94
1	B	62	GYC	CG2-CB2-CA2	-4.87	123.98	129.94
1	A	62	GYC	O2-C2-CA2	-4.61	128.37	130.96
1	B	62	GYC	C2-CA2-N2	-4.36	105.88	108.93
1	D	62	GYC	C2-CA2-N2	-3.90	106.20	108.93
1	D	62	GYC	CG2-CB2-CA2	-3.88	125.19	129.94
1	A	62	GYC	C2-CA2-N2	-3.77	106.29	108.93
1	D	62	GYC	O2-C2-CA2	-3.53	128.98	130.96
1	B	62	GYC	CA2-N2-C1	3.49	108.34	105.77
1	D	62	GYC	CA2-N2-C1	3.47	108.33	105.77
1	C	62	GYC	CG2-CB2-CA2	-3.44	125.73	129.94
1	C	62	GYC	C2-CA2-N2	-3.38	106.57	108.93
1	B	62	GYC	CA3-N3-C1	-3.35	123.14	127.16
1	C	62	GYC	O-C-CA3	-3.33	116.35	126.39
1	C	62	GYC	O2-C2-CA2	-3.08	129.23	130.96
1	A	62	GYC	O-C-CA3	-3.07	117.14	126.39
1	D	62	GYC	O-C-CA3	-2.86	117.74	126.39
1	B	62	GYC	N3-C1-N2	-2.72	109.57	111.45
1	B	62	GYC	O-C-CA3	-2.70	118.24	126.39
1	D	62	GYC	CA3-N3-C1	-2.63	124.01	127.16
1	D	62	GYC	N3-C1-N2	-2.57	109.67	111.45
1	A	62	GYC	CB2-CA2-N2	2.49	132.28	128.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	62	GYC	CA3-N3-C1	-2.32	124.38	127.16
1	A	62	GYC	CA2-N2-C1	2.27	107.45	105.77
1	C	62	GYC	CA1-CB1-SG1	-2.07	109.97	114.44
1	B	62	GYC	CA3-N3-C2	2.05	128.51	123.80

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	62	GYC	C-CA3-N3-C2
1	B	62	GYC	C-CA3-N3-C2
1	D	62	GYC	C-CA3-N3-C2
1	A	62	GYC	C-CA3-N3-C2
1	B	62	GYC	N2-CA2-CB2-CG2
1	A	62	GYC	N2-CA2-CB2-CG2
1	C	62	GYC	N2-CA2-CB2-CG2

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	62	GYC	1	0
1	B	62	GYC	2	0
1	D	62	GYC	1	0
1	A	62	GYC	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 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	GOL	A	225	-	5,5,5	0.34	0	5,5,5	0.42	0
2	GOL	D	227	-	5,5,5	0.44	0	5,5,5	0.32	0
2	GOL	C	225	-	5,5,5	0.31	0	5,5,5	0.31	0
3	SPD	B	226	-	9,9,9	0.28	0	8,8,8	0.77	0
3	SPD	C	226	-	9,9,9	0.29	0	8,8,8	0.67	0
3	SPD	B	225	-	9,9,9	0.28	0	8,8,8	0.93	0
4	EPE	D	225	-	15,15,15	1.03	1 (6%)	18,20,20	2.02	5 (27%)
3	SPD	C	227	-	9,9,9	0.28	0	8,8,8	0.61	0
3	SPD	A	227	-	9,9,9	0.33	0	8,8,8	0.69	0
2	GOL	D	226	-	5,5,5	0.47	0	5,5,5	0.37	0
2	GOL	A	226	-	5,5,5	0.38	0	5,5,5	0.21	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	225	-	-	4/4/4/4	-
2	GOL	D	227	-	-	4/4/4/4	-
2	GOL	C	225	-	-	2/4/4/4	-
3	SPD	B	226	-	-	4/7/7/7	-
3	SPD	C	226	-	-	5/7/7/7	-
3	SPD	B	225	-	-	5/7/7/7	-
4	EPE	D	225	-	-	6/9/19/19	0/1/1/1
3	SPD	C	227	-	-	3/7/7/7	-
3	SPD	A	227	-	-	5/7/7/7	-
2	GOL	D	226	-	-	2/4/4/4	-
2	GOL	A	226	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	225	EPE	C10-S	3.43	1.82	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	225	EPE	C5-N4-C3	4.68	119.36	108.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	225	EPE	C7-N4-C3	3.71	120.72	111.23
4	D	225	EPE	C7-N4-C5	3.39	119.92	111.23
4	D	225	EPE	O2S-S-C10	3.35	110.95	106.92
4	D	225	EPE	C6-N1-C2	2.90	115.35	108.83

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	227	GOL	O1-C1-C2-C3
2	D	227	GOL	C1-C2-C3-O3
2	C	225	GOL	O1-C1-C2-C3
4	D	225	EPE	S-C10-C9-N1
2	D	226	GOL	O1-C1-C2-C3
3	A	227	SPD	N6-C7-C8-C9
3	C	226	SPD	N6-C7-C8-C9
3	B	225	SPD	N6-C7-C8-C9
2	D	227	GOL	O1-C1-C2-O2
2	D	227	GOL	O2-C2-C3-O3
4	D	225	EPE	C9-C10-S-O3S
2	A	225	GOL	O1-C1-C2-C3
2	A	225	GOL	C1-C2-C3-O3
2	A	226	GOL	C1-C2-C3-O3
2	A	225	GOL	O1-C1-C2-O2
2	C	225	GOL	O1-C1-C2-O2
2	D	226	GOL	O1-C1-C2-O2
3	B	226	SPD	C2-C3-C4-C5
2	A	225	GOL	O2-C2-C3-O3
3	C	226	SPD	C2-C3-C4-C5
4	D	225	EPE	C10-C9-N1-C2
2	A	226	GOL	O2-C2-C3-O3
3	A	227	SPD	C8-C7-N6-C5
3	A	227	SPD	C4-C5-N6-C7
3	C	226	SPD	N1-C2-C3-C4
3	C	227	SPD	C7-C8-C9-N10
3	B	226	SPD	C3-C4-C5-N6
4	D	225	EPE	C9-C10-S-O1S
4	D	225	EPE	C9-C10-S-O2S
3	A	227	SPD	N1-C2-C3-C4
3	B	225	SPD	C7-C8-C9-N10
3	A	227	SPD	C7-C8-C9-N10
3	B	225	SPD	C3-C4-C5-N6

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Mol	Chain	Res	Type	Atoms
3	C	227	SPD	C8-C7-N6-C5
3	B	226	SPD	C8-C7-N6-C5
3	C	227	SPD	C4-C5-N6-C7
3	B	226	SPD	C4-C5-N6-C7
3	C	226	SPD	C4-C5-N6-C7
3	B	225	SPD	C4-C5-N6-C7
3	C	226	SPD	C3-C4-C5-N6
3	B	225	SPD	C8-C7-N6-C5
4	D	225	EPE	C8-C7-N4-C3

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	227	GOL	1	0
3	C	226	SPD	1	0
3	B	225	SPD	6	0
3	C	227	SPD	1	0
3	A	227	SPD	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	215/235 (91%)	0.34	16 (7%) 14 14	15, 20, 27, 47	1 (0%)
1	B	214/235 (91%)	0.18	10 (4%) 31 30	14, 19, 26, 40	0
1	C	213/235 (90%)	0.17	8 (3%) 40 40	11, 16, 24, 32	1 (0%)
1	D	224/235 (95%)	0.16	11 (4%) 29 28	11, 16, 25, 36	1 (0%)
All	All	866/940 (92%)	0.21	45 (5%) 27 26	11, 18, 26, 47	3 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	9.3
1	D	0	PRO	6.6
1	B	2	SER	6.4
1	A	2	SER	5.5
1	D	165	GLY	5.4
1	D	-2	THR	5.2
1	D	1	MET	4.8
1	B	218	GLU	4.5
1	A	218	GLU	4.3
1	D	-8	HIS	4.2
1	A	128	ASN	4.0
1	A	67	VAL	3.9
1	A	165	GLY	3.8
1	C	73	GLU	3.8
1	A	73	GLU	3.6
1	D	218	GLU	3.6
1	D	-1	ASP	3.5
1	B	73	GLU	3.4
1	A	68	PHE	3.2
1	B	68	PHE	3.2
1	A	56	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	68	PHE	3.0
1	B	3	VAL	2.7
1	A	202	LYS	2.6
1	D	67	VAL	2.6
1	B	79	PHE	2.5
1	C	2	SER	2.4
1	D	-5	HIS	2.3
1	C	67	VAL	2.3
1	D	79	PHE	2.3
1	B	128	ASN	2.3
1	A	198	ILE	2.2
1	A	3	VAL	2.2
1	B	146	LEU	2.2
1	C	112	ASP	2.2
1	C	3	VAL	2.2
1	A	209[A]	LEU	2.2
1	A	164	GLU	2.1
1	B	56	ILE	2.1
1	C	30	LEU	2.1
1	D	56	ILE	2.1
1	B	67	VAL	2.1
1	A	74	ASN	2.1
1	C	182	VAL	2.0
1	A	5	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	GYC	C	62	21/22	0.94	0.09	15,19,22,27	0
1	GYC	A	62	21/22	0.95	0.08	19,22,27,31	0
1	GYC	D	62	21/22	0.96	0.07	15,17,22,25	0
1	GYC	B	62	21/22	0.96	0.07	17,20,24,29	0

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GOL	C	225	6/6	0.64	0.17	52,55,56,58	0
3	SPD	C	227	10/10	0.72	0.24	39,55,57,57	0
2	GOL	A	226	6/6	0.74	0.21	64,64,64,64	0
2	GOL	D	227	6/6	0.75	0.19	55,58,59,59	0
3	SPD	B	226	10/10	0.78	0.14	50,55,56,56	0
3	SPD	B	225	10/10	0.78	0.16	50,56,60,61	0
3	SPD	A	227	10/10	0.79	0.16	47,51,57,57	0
3	SPD	C	226	10/10	0.79	0.12	59,61,63,63	0
2	GOL	D	226	6/6	0.80	0.18	59,61,62,63	0
2	GOL	A	225	6/6	0.81	0.19	79,79,79,80	0
4	EPE	D	225	15/15	0.89	0.36	55,67,69,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.