



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 1, 2021 – 06:05 PM EDT

PDB ID : 7LCE
Title : Structure of D-Glucosamine-6-phosphate Ammonia-lyase
Authors : Phillips, R.S.
Deposited on : 2021-01-10
Resolution : 2.58 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 : 2.19
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) : 2.19

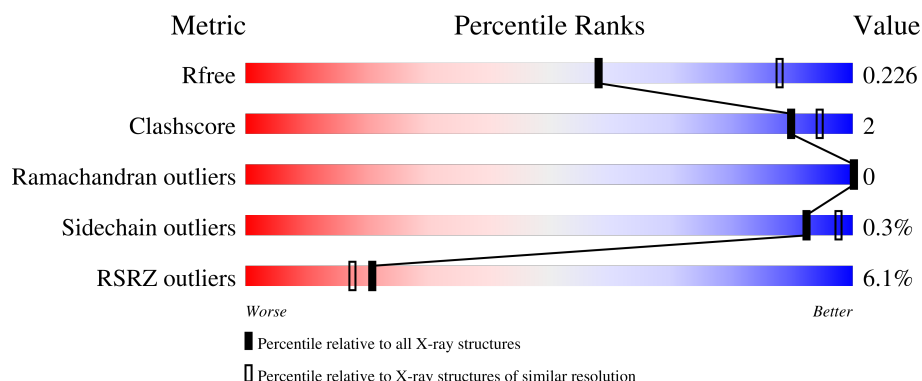
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 of 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	369	<div> <div>0%</div> <div>95%</div> <div>5%</div> </div>
1	B	369	<div> <div>4%</div> <div>97%</div> <div>.</div> </div>
1	C	369	<div> <div>12%</div> <div>95%</div> <div>5%</div> </div>
1	D	369	<div> <div>8%</div> <div>93%</div> <div>7%</div> </div>

2 Entry composition [i](#)

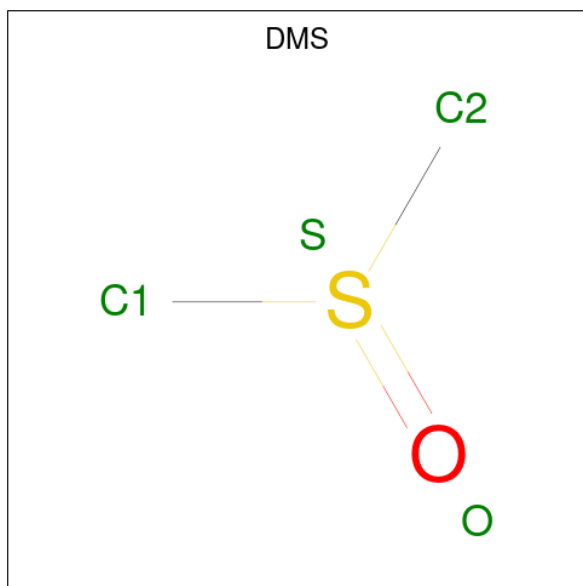
There are 4 unique types of molecules in this entry. The entry contains 11310 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-glucosaminatase-6-phosphate ammonia lyase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	P	S	0	1	0
			2782	1736	500	530	1	15			
1	B	369	Total	C	N	O	P	S	0	0	0
			2774	1731	497	530	1	15			
1	C	369	Total	C	N	O	P	S	0	1	0
			2784	1737	500	531	1	15			
1	D	369	Total	C	N	O	P	S	0	1	0
			2780	1735	498	531	1	15			

- Molecule 2 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 4	C 2	O 1	S 1	0	0
2	B	1	Total 4	C 2	O 1	S 1	0	0
2	C	1	Total 4	C 2	O 1	S 1	0	0
2	D	1	Total 4	C 2	O 1	S 1	0	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total 6	Ca 6	0	0
3	B	4	Total 4	Ca 4	0	0
3	C	3	Total 3	Ca 3	0	0
3	D	3	Total 3	Ca 3	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total 41	O 41	0	0
4	B	34	Total 34	O 34	0	0
4	C	26	Total 26	O 26	0	0
4	D	49	Total 49	O 49	0	0

3 Residue-property plots [i](#)

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

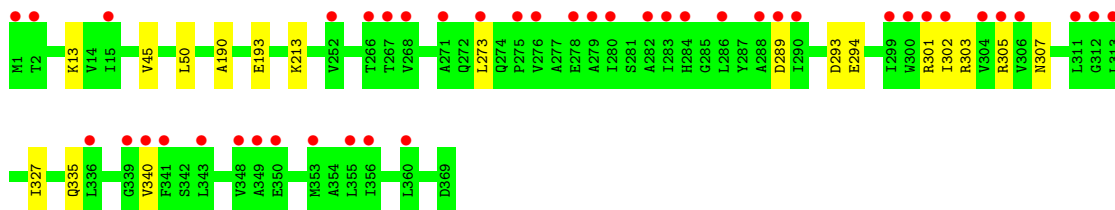
- Molecule 1: D-glucosamine-6-phosphate ammonia lyase



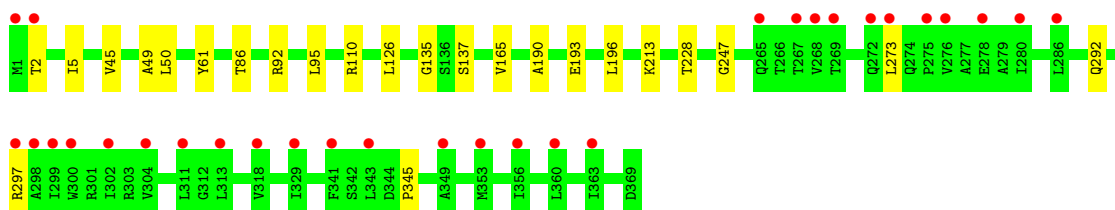
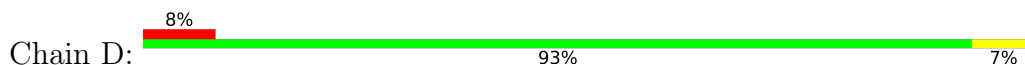
- Molecule 1: D-glucosamine-6-phosphate ammonia lyase



- Molecule 1: D-glucosamine-6-phosphate ammonia lyase



- Molecule 1: D-glucosamine-6-phosphate ammonia lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	60.37Å 171.57Å 195.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.37 – 2.58 60.37 – 2.58	Depositor EDS
% Data completeness (in resolution range)	75.7 (60.37-2.58) 75.7 (60.37-2.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.193 , 0.229 0.192 , 0.226	Depositor DCC
R_{free} test set	2000 reflections (4.09%)	wwPDB-VP
Wilson B-factor (Å ²)	78.6	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11310	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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: LLP, CA, DMS

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	0.24	0/2801	0.50	0/3794
1	B	0.25	0/2790	0.49	0/3780
1	C	0.25	0/2801	0.49	0/3795
1	D	0.25	0/2799	0.49	0/3792
All	All	0.25	0/11191	0.49	0/15161

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	2782	0	2791	13	0
1	B	2774	0	2778	7	0
1	C	2784	0	2784	10	0
1	D	2780	0	2786	14	0
2	A	12	0	18	0	0
2	B	4	0	6	0	0
2	C	4	0	6	0	0
2	D	4	0	6	0	0
3	A	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	0	0	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
4	A	41	0	0	0	0
4	B	34	0	0	3	0
4	C	26	0	0	0	0
4	D	49	0	0	0	0
All	All	11310	0	11175	39	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 2.

All (39) 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:34:MET:SD	4:B:502:HOH:O	2.40	0.79
1:D:86:THR:HA	1:D:92:ARG:HD3	1.86	0.57
1:A:273:LEU:HD22	1:A:302:ILE:HD11	1.87	0.56
1:C:305:ARG:NH1	1:C:307:ASN:HB2	2.21	0.55
1:A:280:ILE:HA	1:A:283:ILE:HD12	1.92	0.52
1:C:13:LYS:HD3	1:C:327:ILE:HD12	1.92	0.52
1:D:61:TYR:HB3	1:D:196:LEU:HD13	1.92	0.52
1:A:256:TYR:CE1	1:D:5:ILE:HD11	2.46	0.50
1:C:273:LEU:HD22	1:C:302:ILE:HD11	1.94	0.50
1:A:97:PRO:HG3	1:D:126:LEU:HD23	1.93	0.50
1:A:126:LEU:HD22	1:D:126:LEU:HD22	1.95	0.49
1:A:45:VAL:HG11	1:A:50:LEU:HD22	1.95	0.48
1:B:15:ILE:HG12	4:B:505:HOH:O	2.13	0.48
1:C:45:VAL:HG11	1:C:50:LEU:HD22	1.96	0.47
1:C:289:ASP:OD2	1:C:303:ARG:NH2	2.47	0.47
1:A:79:ILE:HG23	1:A:236:GLN:HG3	1.97	0.46
1:C:190:ALA:HB1	1:C:193:GLU:HB2	1.98	0.46
1:D:110:ARG:NH1	1:D:135:GLY:O	2.45	0.45
1:A:289:ASP:OD2	1:A:303:ARG:NH2	2.44	0.45
1:D:45:VAL:HG11	1:D:50:LEU:HD22	1.98	0.45
1:A:18:CYS:HA	1:A:332:ARG:HG3	1.99	0.45
1:D:273:LEU:HD11	1:D:345:PRO:HG3	1.98	0.45
1:C:305:ARG:HH12	1:C:307:ASN:HB2	1.80	0.45
1:B:301:ARG:HH21	1:B:346:ARG:NH2	2.15	0.44
1:B:274:GLN:HA	1:B:290:ILE:HD11	2.00	0.44
1:A:205:ASP:O	1:A:225:GLY:HA3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:LEU:HD23	1:D:95:LEU:HA	1.86	0.43
1:A:97:PRO:HB3	1:A:126:LEU:O	2.19	0.43
1:C:293:ASP:OD1	1:C:294:GLU:N	2.52	0.42
1:B:298:ALA:HA	4:B:509:HOH:O	2.19	0.42
1:A:11:LEU:HD22	1:D:49:ALA:HB1	2.02	0.42
1:C:335:GLN:HB2	1:C:340:VAL:HG12	2.01	0.42
1:A:219:THR:OG1	1:D:247:GLY:HA3	2.20	0.41
1:D:137:SER:O	1:D:165:VAL:HG21	2.20	0.41
1:C:293:ASP:CG	1:C:301:ARG:HH12	2.23	0.41
1:D:190:ALA:HB1	1:D:193:GLU:HB2	2.03	0.41
1:B:15:ILE:HD12	1:B:329:ILE:HG12	2.01	0.41
1:B:293:ASP:HB3	1:B:296:GLY:O	2.21	0.41
1:D:292:GLN:OE1	1:D:297:ARG:HG2	2.22	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	367/369 (100%)	357 (97%)	10 (3%)	0	100	100
1	B	366/369 (99%)	359 (98%)	7 (2%)	0	100	100
1	C	367/369 (100%)	360 (98%)	7 (2%)	0	100	100
1	D	367/369 (100%)	357 (97%)	10 (3%)	0	100	100
All	All	1467/1476 (99%)	1433 (98%)	34 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	278/277 (100%)	278 (100%)	0	100	100
1	B	277/277 (100%)	276 (100%)	1 (0%)	91	97
1	C	278/277 (100%)	278 (100%)	0	100	100
1	D	278/277 (100%)	276 (99%)	2 (1%)	84	93
All	All	1111/1108 (100%)	1108 (100%)	3 (0%)	92	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	366	HIS
1	D	2	THR
1	D	228	THR

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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	LLP	B	213	1	23,24,25	2.62	6 (26%)	25,32,34	1.28	3 (12%)
1	LLP	C	213	1	23,24,25	2.60	7 (30%)	25,32,34	1.43	3 (12%)
1	LLP	A	213	1	23,24,25	2.56	6 (26%)	25,32,34	1.48	5 (20%)
1	LLP	D	213	1	23,24,25	2.61	6 (26%)	25,32,34	1.27	3 (12%)

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	LLP	B	213	1	-	4/16/17/19	0/1/1/1
1	LLP	C	213	1	-	5/16/17/19	0/1/1/1
1	LLP	A	213	1	-	4/16/17/19	0/1/1/1
1	LLP	D	213	1	-	4/16/17/19	0/1/1/1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	213	LLP	C4-C4'	8.19	1.62	1.46
1	D	213	LLP	C4-C4'	8.08	1.62	1.46
1	C	213	LLP	C4-C4'	7.85	1.61	1.46
1	A	213	LLP	C4-C4'	7.84	1.61	1.46
1	B	213	LLP	C4'-NZ	4.94	1.43	1.27
1	D	213	LLP	C4'-NZ	4.88	1.43	1.27
1	C	213	LLP	C4'-NZ	4.77	1.43	1.27
1	A	213	LLP	C4'-NZ	4.76	1.43	1.27
1	C	213	LLP	C4-C5	-4.57	1.36	1.42
1	B	213	LLP	C4-C5	-4.07	1.36	1.42
1	D	213	LLP	C4-C5	-4.02	1.36	1.42
1	A	213	LLP	C4-C5	-4.01	1.36	1.42
1	C	213	LLP	C2'-C2	3.50	1.56	1.50
1	A	213	LLP	C2'-C2	3.46	1.56	1.50
1	D	213	LLP	C2'-C2	3.46	1.56	1.50
1	B	213	LLP	C2'-C2	3.37	1.56	1.50
1	A	213	LLP	C6-N1	2.99	1.40	1.34
1	C	213	LLP	C6-N1	2.95	1.40	1.34
1	D	213	LLP	C6-N1	2.92	1.40	1.34
1	B	213	LLP	C6-N1	2.89	1.40	1.34
1	A	213	LLP	C5'-C5	2.15	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	213	LLP	C5'-C5	2.11	1.56	1.50
1	B	213	LLP	C5'-C5	2.09	1.56	1.50
1	C	213	LLP	CB-CA	-2.07	1.50	1.53
1	C	213	LLP	C5'-C5	2.01	1.56	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	213	LLP	CE-NZ-C4'	-3.67	107.62	118.90
1	A	213	LLP	C4-C4'-NZ	-3.59	107.84	124.31
1	C	213	LLP	C4-C4'-NZ	-3.44	108.52	124.31
1	B	213	LLP	C4-C4'-NZ	-3.34	108.99	124.31
1	D	213	LLP	C4-C4'-NZ	-3.05	110.28	124.31
1	C	213	LLP	C3-C4-C5	2.89	120.47	118.26
1	C	213	LLP	C5-C6-N1	-2.71	119.31	123.82
1	B	213	LLP	CE-NZ-C4'	-2.58	110.97	118.90
1	D	213	LLP	CE-NZ-C4'	-2.57	111.02	118.90
1	A	213	LLP	C3-C4-C5	2.36	120.08	118.26
1	D	213	LLP	C5-C6-N1	-2.36	119.89	123.82
1	B	213	LLP	C5-C6-N1	-2.34	119.91	123.82
1	A	213	LLP	C5-C6-N1	-2.33	119.94	123.82
1	A	213	LLP	CD-CG-CB	-2.09	106.23	113.62

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	213	LLP	C-CA-CB-CG
1	B	213	LLP	C-CA-CB-CG
1	C	213	LLP	C-CA-CB-CG
1	D	213	LLP	C-CA-CB-CG
1	A	213	LLP	C4-C4'-NZ-CE
1	B	213	LLP	C4-C4'-NZ-CE
1	D	213	LLP	C4-C4'-NZ-CE
1	C	213	LLP	C4-C4'-NZ-CE
1	A	213	LLP	N-CA-CB-CG
1	D	213	LLP	C5'-OP4-P-OP1
1	A	213	LLP	C3-C4-C4'-NZ
1	B	213	LLP	C3-C4-C4'-NZ
1	C	213	LLP	C3-C4-C4'-NZ
1	D	213	LLP	C3-C4-C4'-NZ
1	C	213	LLP	CD-CE-NZ-C4'

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Mol	Chain	Res	Type	Atoms
1	B	213	LLP	N-CA-CB-CG
1	C	213	LLP	N-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 16 are monoatomic - leaving 6 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	$\# Z > 2$
2	DMS	A	401	-	3,3,3	0.67	0	3,3,3	0.51	0
2	DMS	A	402	-	3,3,3	0.67	0	3,3,3	0.55	0
2	DMS	A	403	-	3,3,3	0.67	0	3,3,3	0.57	0
2	DMS	C	401	-	3,3,3	0.68	0	3,3,3	0.57	0
2	DMS	B	401	-	3,3,3	0.68	0	3,3,3	0.54	0
2	DMS	D	401	-	3,3,3	0.67	0	3,3,3	0.56	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	368/369 (99%)	0.09	3 (0%) 86 85	47, 72, 111, 167	0
1	B	368/369 (99%)	0.30	14 (3%) 40 36	46, 80, 135, 176	0
1	C	368/369 (99%)	0.59	43 (11%) 4 3	49, 93, 160, 190	0
1	D	368/369 (99%)	0.30	30 (8%) 11 9	47, 75, 133, 169	0
All	All	1472/1476 (99%)	0.32	90 (6%) 21 17	46, 78, 147, 190	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	313	LEU	7.6
1	C	356	ILE	6.7
1	C	283	ILE	5.8
1	D	298	ALA	5.8
1	C	306	VAL	5.7
1	C	302	ILE	5.7
1	B	1	MET	5.6
1	C	1	MET	5.4
1	C	280	ILE	5.3
1	D	1	MET	5.1
1	D	343	LEU	5.0
1	B	280	ILE	5.0
1	D	286	LEU	4.7
1	C	304	VAL	4.6
1	C	360	LEU	4.5
1	B	283	ILE	4.4
1	C	15	ILE	4.4
1	D	280	ILE	4.4
1	D	341	PHE	4.3
1	C	273	LEU	4.3
1	D	356	ILE	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	312	GLY	4.1
1	C	311	LEU	4.1
1	D	273	LEU	4.1
1	C	268	VAL	4.0
1	D	267	THR	3.9
1	C	286	LEU	3.9
1	C	305	ARG	3.9
1	D	311	LEU	3.9
1	C	288	ALA	3.8
1	D	313	LEU	3.7
1	C	341	PHE	3.7
1	D	276	VAL	3.7
1	C	266	THR	3.6
1	C	267	THR	3.6
1	D	304	VAL	3.6
1	C	289	ASP	3.5
1	D	268	VAL	3.4
1	C	339	GLY	3.4
1	C	276	VAL	3.4
1	C	343	LEU	3.3
1	C	284	HIS	3.3
1	C	355	LEU	3.2
1	A	262	HIS	3.2
1	C	300	TRP	3.2
1	B	360	LEU	3.2
1	D	272	GLN	3.2
1	C	353	MET	3.1
1	D	265	GLN	3.1
1	D	360	LEU	3.0
1	C	279	ALA	3.0
1	B	276	VAL	3.0
1	D	363	ILE	2.9
1	D	278	GLU	2.9
1	C	349	ALA	2.9
1	B	286	LEU	2.8
1	C	299	ILE	2.8
1	C	282	ALA	2.8
1	D	297	ARG	2.8
1	C	336	LEU	2.7
1	C	340	VAL	2.6
1	D	269	THR	2.6
1	C	301	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	256	TYR	2.6
1	D	318	VAL	2.6
1	C	271	ALA	2.5
1	C	348	VAL	2.4
1	A	299	ILE	2.4
1	D	302	ILE	2.4
1	B	268	VAL	2.3
1	D	275	PRO	2.3
1	B	313	LEU	2.3
1	B	322	LEU	2.3
1	D	353	MET	2.2
1	D	329	ILE	2.2
1	B	341	PHE	2.2
1	C	350	GLU	2.2
1	C	2	THR	2.2
1	B	311	LEU	2.2
1	C	275	PRO	2.2
1	C	290	ILE	2.2
1	B	299	ILE	2.2
1	D	299	ILE	2.2
1	B	285	GLY	2.1
1	C	278	GLU	2.1
1	D	2	THR	2.1
1	D	300	TRP	2.1
1	C	252	VAL	2.1
1	D	349	ALA	2.1
1	B	300	TRP	2.0

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

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	LLP	A	213	24/25	0.96	0.16	65,81,102,102	0
1	LLP	B	213	24/25	0.96	0.17	58,77,90,105	0
1	LLP	D	213	24/25	0.96	0.17	61,74,82,85	0
1	LLP	C	213	24/25	0.97	0.17	77,93,104,113	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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	DMS	B	401	4/4	0.62	0.28	75,85,109,149	0
2	DMS	A	403	4/4	0.75	0.33	110,114,125,153	0
2	DMS	A	402	4/4	0.76	0.33	86,94,100,131	0
3	CA	B	404	1/1	0.80	0.23	115,115,115,115	0
3	CA	B	402	1/1	0.84	0.09	141,141,141,141	0
2	DMS	C	401	4/4	0.84	0.70	81,83,85,130	0
2	DMS	D	401	4/4	0.88	0.17	83,87,113,136	0
2	DMS	A	401	4/4	0.90	0.26	107,114,115,138	0
3	CA	B	405	1/1	0.90	0.07	102,102,102,102	0
3	CA	A	407	1/1	0.91	0.07	108,108,108,108	0
3	CA	C	404	1/1	0.91	0.08	110,110,110,110	0
3	CA	A	409	1/1	0.92	0.06	97,97,97,97	0
3	CA	D	403	1/1	0.92	0.11	89,89,89,89	0
3	CA	A	406	1/1	0.93	0.19	94,94,94,94	0
3	CA	B	403	1/1	0.93	0.10	86,86,86,86	0
3	CA	D	402	1/1	0.94	0.09	89,89,89,89	0
3	CA	A	404	1/1	0.94	0.13	87,87,87,87	0
3	CA	D	404	1/1	0.95	0.10	93,93,93,93	0
3	CA	A	408	1/1	0.96	0.02	125,125,125,125	1
3	CA	C	402	1/1	0.97	0.13	91,91,91,91	0
3	CA	C	403	1/1	0.97	0.15	102,102,102,102	0
3	CA	A	405	1/1	0.97	0.14	79,79,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.