



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:11 PM GMT

PDB ID : 4G36
Title : Photinus pyralis luciferase in the adenylate-forming conformation bound to DLSA
Authors : Sundlov, J.A.; Branchini, B.R.; Gulick, A.M.
Deposited on : 2012-07-13
Resolution : 2.62 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>

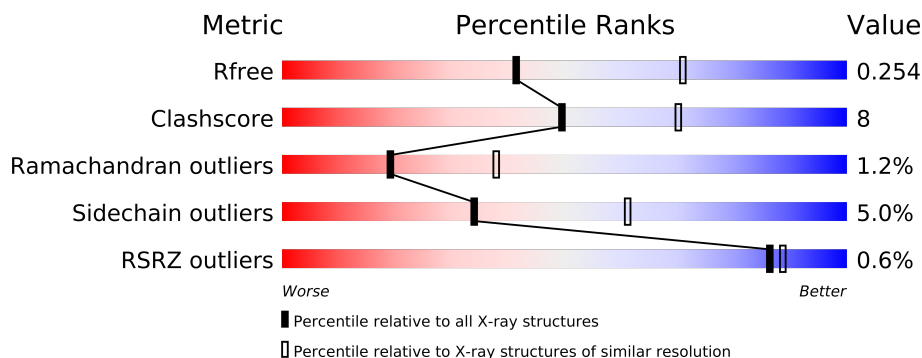
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
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 2.62 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}	66092	1940 (2.64-2.60)
Clashscore	79885	2404 (2.64-2.60)
Ramachandran outliers	78287	2360 (2.64-2.60)
Sidechain outliers	78261	2360 (2.64-2.60)
RSRZ outliers	66119	1939 (2.64-2.60)

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.

Mol	Chain	Length	Quality of chain
1	A	555	
1	B	555	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8222 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.

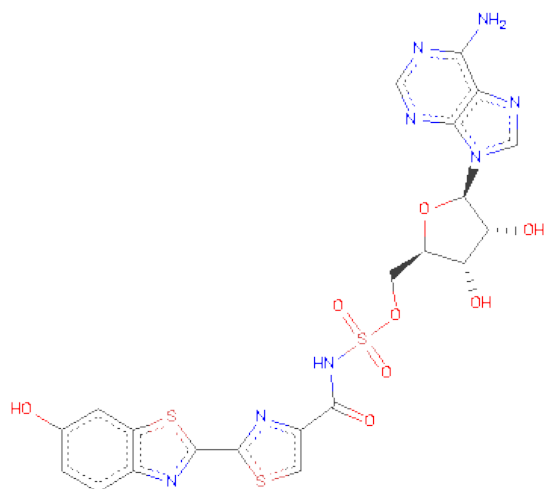
- Molecule 1 is a protein called Luciferin 4-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	522	Total	C	N	O	S	0	0	0
			3942	2541	651	734	16			
1	B	543	Total	C	N	O	S	0	0	0
			4128	2658	687	766	17			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP P08659
A	-3	PRO	-	EXPRESSION TAG	UNP P08659
A	-2	LEU	-	EXPRESSION TAG	UNP P08659
A	-1	GLY	-	EXPRESSION TAG	UNP P08659
A	0	SER	-	EXPRESSION TAG	UNP P08659
B	-4	GLY	-	EXPRESSION TAG	UNP P08659
B	-3	PRO	-	EXPRESSION TAG	UNP P08659
B	-2	LEU	-	EXPRESSION TAG	UNP P08659
B	-1	GLY	-	EXPRESSION TAG	UNP P08659
B	0	SER	-	EXPRESSION TAG	UNP P08659

- Molecule 2 is 5'-O-[N-(DEHYDROLUCIFERYL)-SULFAMOYL]ADENOSINE (three-letter code: SLU) (formula: C₂₁H₁₈N₈O₈S₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			40	21	8	8	3		
2	B	1	Total	C	N	O	S	0	0
			40	21	8	8	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	46	Total	O	0	0
			46	46		
3	B	26	Total	O	0	0
			26	26		

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	93.04Å 93.04Å 297.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	54.66 – 2.62 54.66 – 2.62	Depositor EDS
% Data completeness (in resolution range)	99.4 (54.66-2.62) 99.4 (54.66-2.62)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.42 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_837)	Depositor
R, R_{free}	0.187 , 0.238 0.202 , 0.254	Depositor DCC
R_{free} test set	2301 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 25.7	EDS
Estimated twinning fraction	0.079 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	1 of 45569 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8222	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SLU

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	0.41	0/4028	0.59	0/5478
1	B	0.39	0/4218	0.57	0/5731
All	All	0.40	0/8246	0.58	0/11209

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	501	TYR	Sidechain

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	3942	0	3837	73	0
1	B	4128	0	4071	49	0
2	A	40	0	18	1	0
2	B	40	0	18	1	0
3	A	46	0	0	3	0
3	B	26	0	0	1	0
All	All	8222	0	7944	122	0

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

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:495:GLU:HG3	1:A:516:VAL:HB	1.59	0.84
1:A:275:ARG:NH1	3:A:718:HOH:O	2.13	0.80
1:A:238:LEU:HB2	1:A:282:ILE:HD13	1.64	0.79
1:A:79:VAL:HG13	1:A:123:PRO:HB3	1.70	0.73
1:A:107:ASP:OD1	3:A:732:HOH:O	2.14	0.65
1:B:98:ILE:HG13	1:B:100:VAL:HG23	1.78	0.65
1:A:44:ASP:HB2	1:A:86:LEU:HD21	1.81	0.62
1:B:77:ARG:NH1	1:B:122:GLN:O	2.32	0.62
1:A:295:PHE:HA	1:A:298:SER:HB2	1.82	0.60
1:A:338:GLN:HG2	1:A:362:VAL:HA	1.82	0.60
1:A:524:LYS:HG2	1:A:530:LEU:HD23	1.82	0.59
1:B:81:CYS:HB2	1:B:105:ALA:HB3	1.84	0.59
1:A:62:ARG:HG2	1:A:168:VAL:HG11	1.85	0.58
1:A:4:ALA:HB1	1:A:6:ASN:HB2	1.85	0.58
1:B:79:VAL:HG21	1:B:118:MET:HE3	1.87	0.57
1:A:454:LEU:HD12	1:A:469:VAL:HG11	1.86	0.56
1:B:524:LYS:HG2	1:B:530:LEU:HD23	1.88	0.56
1:A:58:GLU:O	1:A:62:ARG:HB2	2.05	0.56
1:A:241:VAL:HG12	1:A:288:VAL:HG13	1.88	0.55
1:B:79:VAL:HG23	1:B:123:PRO:HB3	1.89	0.55
1:A:43:THR:HB	1:A:263:VAL:HG22	1.89	0.55
1:A:290:THR:HG23	1:A:528:GLY:HA2	1.90	0.54
2:A:601:SLU:H212	2:A:601:SLU:H36	1.89	0.54
1:A:62:ARG:HG3	1:A:172:LEU:HD11	1.90	0.54
1:B:114:LEU:HD23	1:B:135:LYS:HB2	1.89	0.54
1:B:115:LEU:HD13	1:B:139:VAL:HG13	1.90	0.54
1:B:111:GLU:HG2	1:B:135:LYS:HB3	1.91	0.53
1:A:340:TYR:HB3	1:A:350:LEU:HB2	1.92	0.52
1:A:4:ALA:HB1	1:A:6:ASN:H	1.75	0.52
1:A:51:ILE:HG12	1:A:56:TYR:HD2	1.74	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:302:ASP:OD2	1:A:330:ARG:NH2	2.43	0.51
1:A:338:GLN:HG3	1:A:339:GLY:N	2.25	0.51
1:A:187:ASP:HB3	1:A:190:LYS:HB2	1.93	0.51
1:A:172:LEU:HD22	1:A:173:PRO:HD2	1.93	0.50
1:B:322:GLU:HG2	1:B:460:GLN:HE21	1.76	0.50
1:B:313:ALA:HA	1:B:337:ARG:O	2.12	0.49
1:A:292:PHE:HD2	1:A:327:VAL:HG21	1.76	0.49
1:A:225:PRO:HB3	1:A:230:GLN:HB2	1.94	0.49
1:A:65:GLU:HG2	1:A:176:PHE:CE1	2.48	0.49
1:A:498:ILE:O	1:A:502:VAL:N	2.36	0.49
1:A:106:ASN:HB3	1:A:109:TYR:CG	2.49	0.48
1:B:133:LEU:HD13	1:B:151:ILE:HD11	1.93	0.48
1:A:62:ARG:HG3	1:A:172:LEU:CD1	2.44	0.48
1:A:344:GLU:N	1:A:344:GLU:OE1	2.45	0.48
1:B:455:GLU:HG3	1:B:469:VAL:HG23	1.94	0.48
1:A:443:LYS:HE2	1:A:448:GLN:HB2	1.95	0.48
1:A:454:LEU:HA	1:A:457:ILE:HD12	1.97	0.47
1:B:110:ASN:ND2	1:B:113:GLU:OE1	2.47	0.47
1:B:484:VAL:HG13	1:B:519:VAL:HG21	1.95	0.47
1:A:272:LEU:HD11	1:A:275:ARG:HH21	1.77	0.47
1:A:290:THR:CG2	1:A:528:GLY:HA2	2.44	0.47
1:A:237:ILE:HD12	1:A:284:SER:HB3	1.97	0.46
1:A:517:VAL:HB	1:A:539:LEU:HD23	1.96	0.46
1:B:195:ILE:HA	1:B:208:VAL:O	2.16	0.46
1:A:311:GLU:HA	1:A:333:LEU:HD22	1.97	0.46
1:B:194:LEU:HD13	1:B:250:PHE:CZ	2.51	0.46
1:B:531:ASP:OD2	1:B:534:LYS:HD3	2.15	0.46
1:B:442:ILE:HG23	1:B:479:GLU:HG2	1.97	0.46
1:A:338:GLN:HE21	1:A:338:GLN:HB2	1.50	0.46
1:B:74:THR:HG22	1:B:99:GLY:O	2.15	0.46
1:A:67:MET:HB3	1:A:72:LEU:HD12	1.96	0.46
1:A:358:LYS:HB2	1:A:358:LYS:HE3	1.73	0.46
1:A:321:LYS:HD2	1:A:359:PRO:HB2	1.97	0.46
1:B:65:GLU:HG2	1:B:176:PHE:CE1	2.51	0.45
1:B:465:PHE:CD2	1:B:466:ASP:HB2	2.51	0.45
1:A:160:GLY:O	3:A:707:HOH:O	2.21	0.45
1:B:293:SER:OG	1:B:294:PHE:N	2.49	0.45
1:B:77:ARG:N	1:B:77:ARG:HD3	2.31	0.45
1:B:527:THR:OG1	1:B:529:LYS:HE2	2.17	0.45
1:B:46:HIS:CD2	1:B:272:LEU:HD21	2.52	0.45
1:A:76:HIS:C	1:A:77:ARG:HD3	2.37	0.45
1:B:372:LYS:HE2	1:B:417:TRP:CE2	2.52	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:77:ARG:N	1:A:77:ARG:HD3	2.32	0.44
1:A:511:LYS:O	1:A:513:ARG:N	2.46	0.44
1:A:51:ILE:HG22	1:A:86:LEU:HD13	2.00	0.44
1:B:480:LEU:HA	1:B:481:PRO:HD3	1.82	0.44
1:A:79:VAL:HG22	1:A:126:VAL:HG22	1.98	0.44
1:B:257:ILE:HD13	1:B:257:ILE:HA	1.81	0.44
1:B:113:GLU:HG2	1:B:113:GLU:H	1.68	0.44
1:B:470:ALA:HB3	1:B:539:LEU:HD12	1.99	0.44
1:B:296:ALA:HA	1:B:327:VAL:HG22	1.99	0.44
1:B:471:GLY:HA3	1:B:479:GLU:HG2	2.00	0.43
1:A:272:LEU:HD12	1:A:275:ARG:HE	1.84	0.43
1:A:512:LEU:O	1:A:514:GLY:N	2.50	0.43
1:B:455:GLU:O	1:B:459:LEU:HG	2.19	0.43
1:A:443:LYS:HD3	1:A:443:LYS:HA	1.86	0.43
1:A:274:LEU:HD13	1:A:300:LEU:HB2	2.01	0.43
1:A:247:PHE:CE1	1:A:286:LEU:HD11	2.54	0.43
1:A:524:LYS:HE2	1:A:530:LEU:HD21	1.99	0.42
1:A:128:VAL:HG22	1:A:129:SER:O	2.19	0.42
1:A:247:PHE:HE1	1:A:286:LEU:HD11	1.84	0.42
1:A:512:LEU:C	1:A:514:GLY:H	2.22	0.42
1:A:499:VAL:HG22	1:A:512:LEU:HD23	2.02	0.42
1:B:488:GLU:O	1:B:490:GLY:N	2.45	0.42
1:A:86:LEU:HD23	1:A:264:LEU:HD21	2.00	0.42
1:A:225:PRO:HB2	1:B:225:PRO:HG3	2.02	0.42
1:B:450:ALA:HA	1:B:451:PRO:HD2	1.80	0.42
1:A:77:ARG:HG2	1:A:121:SER:O	2.20	0.42
1:B:277:LEU:HA	1:B:282:ILE:HB	2.01	0.42
1:B:370:GLU:OE1	1:B:393:ARG:NH2	2.35	0.42
1:A:238:LEU:HD22	1:A:273:PHE:CE2	2.55	0.41
1:B:77:ARG:NH2	1:B:122:GLN:HB2	2.35	0.41
1:B:218:ARG:HG2	1:B:250:PHE:HB3	2.02	0.41
1:A:300:LEU:HB3	1:A:304:TYR:HE2	1.85	0.41
1:A:88:PHE:O	1:A:91:PRO:HD2	2.21	0.41
1:B:343:THR:N	2:B:601:SLU:H40	2.19	0.41
1:A:130:LYS:HD2	1:A:158:TYR:CE1	2.55	0.41
1:A:192:ILE:HA	1:A:211:PRO:HA	2.03	0.41
1:A:15:TYR:HA	1:A:16:PRO:HD3	1.81	0.41
1:A:298:SER:O	1:A:330:ARG:NH1	2.53	0.41
1:A:501:TYR:O	1:A:505:GLN:HG2	2.20	0.41
1:A:51:ILE:HG12	1:A:56:TYR:CD2	2.54	0.41
1:B:299:THR:HA	1:B:330:ARG:NH2	2.36	0.41
1:B:302:ASP:OD2	1:B:330:ARG:NH2	2.53	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:32:ARG:NH2	3:B:714:HOH:O	2.54	0.41
1:A:115:LEU:O	1:A:119:ASN:HB2	2.20	0.41
1:A:257:ILE:HD13	1:A:257:ILE:HA	1.96	0.41
1:B:21:THR:O	1:B:25:GLN:HG3	2.22	0.40
1:B:44:ASP:HB3	1:B:47:ILE:HG12	2.02	0.40
1:A:358:LYS:HA	1:A:359:PRO:HD2	1.90	0.40
1:B:287:LEU:O	1:B:314:SER:HA	2.22	0.40
1:B:20:GLY:O	1:B:188:ARG:HD3	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	514/555 (93%)	474 (92%)	32 (6%)	8 (2%)	14	27
1	B	539/555 (97%)	500 (93%)	34 (6%)	5 (1%)	25	48
All	All	1053/1110 (95%)	974 (92%)	66 (6%)	13 (1%)	19	38

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	ALA
1	B	489	HIS
1	B	547	LYS
1	A	301	ILE
1	A	513	ARG
1	A	107	ASP
1	A	346	THR
1	A	435	VAL
1	A	481	PRO
1	B	476	ASP
1	B	477	ALA
1	A	92	VAL

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Mol	Chain	Res	Type
1	B	490	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	408/466 (88%)	390 (96%)	18 (4%)	39	68
1	B	432/466 (93%)	408 (94%)	24 (6%)	30	54
All	All	840/932 (90%)	798 (95%)	42 (5%)	34	61

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	ILE
1	A	56	TYR
1	A	77	ARG
1	A	109	TYR
1	A	170	SER
1	A	172	LEU
1	A	189	ASP
1	A	247	PHE
1	A	299	THR
1	A	338	GLN
1	A	349	ILE
1	A	357	ASP
1	A	387	ARG
1	A	399	SER
1	A	436	ASP
1	A	445	LYS
1	A	474	ASP
1	A	508	THR
1	B	62	ARG
1	B	77	ARG
1	B	109	TYR
1	B	135	LYS
1	B	140	GLN
1	B	154	SER

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Mol	Chain	Res	Type
1	B	155	LYS
1	B	166	THR
1	B	194	LEU
1	B	247	PHE
1	B	257	ILE
1	B	283	GLN
1	B	308	ASN
1	B	310	HIS
1	B	320	SER
1	B	321	LYS
1	B	349	ILE
1	B	357	ASP
1	B	362	VAL
1	B	474	ASP
1	B	502	VAL
1	B	508	THR
1	B	512	LEU
1	B	543	LYS

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 ⓘ

2 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	SLU	A	601	-	45,45,45	2.42	16 (35%)	67,68,68	4.29	28 (41%)
2	SLU	B	601	-	45,45,45	2.41	16 (35%)	67,68,68	4.08	33 (49%)

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	SLU	A	601	-	-	1/17/39/39	0/2/6/6
2	SLU	B	601	-	-	1/17/39/39	0/2/6/6

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	SLU	C16-N40	7.36	1.47	1.39
2	B	601	SLU	C16-N40	6.62	1.46	1.39
2	A	601	SLU	O20-S17	-6.19	1.51	1.58
2	B	601	SLU	O20-S17	-6.13	1.51	1.58
2	B	601	SLU	C8-C11	-5.83	1.38	1.45
2	A	601	SLU	C8-C11	-5.60	1.38	1.45
2	A	601	SLU	S17-N40	4.70	1.65	1.60
2	B	601	SLU	S17-N40	4.44	1.65	1.60
2	B	601	SLU	C24-C25	-4.10	1.47	1.53
2	A	601	SLU	C24-C25	-3.81	1.48	1.53
2	B	601	SLU	C14-N15	-3.69	1.30	1.38
2	B	601	SLU	O19-S17	3.48	1.45	1.42
2	B	601	SLU	C34-N38	3.30	1.45	1.35
2	A	601	SLU	C14-N15	-3.23	1.31	1.38
2	A	601	SLU	C34-N38	3.21	1.45	1.35
2	A	601	SLU	C4-S9	3.08	1.77	1.74
2	B	601	SLU	C4-S9	3.05	1.76	1.74
2	B	601	SLU	O18-S17	3.03	1.45	1.42
2	A	601	SLU	C13-S12	3.03	1.75	1.70
2	A	601	SLU	O19-S17	2.93	1.45	1.42
2	B	601	SLU	C13-S12	2.58	1.74	1.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	SLU	C5-C4	2.55	1.41	1.37
2	B	601	SLU	O26-C22	-2.36	1.39	1.45
2	B	601	SLU	C24-C23	-2.34	1.46	1.53
2	A	601	SLU	C24-C23	-2.29	1.47	1.53
2	A	601	SLU	C3-N7	-2.19	1.31	1.38
2	B	601	SLU	C5-C4	2.17	1.41	1.37
2	B	601	SLU	C3-N7	-2.14	1.31	1.38
2	B	601	SLU	O28-C24	-2.13	1.37	1.43
2	A	601	SLU	O18-S17	2.12	1.44	1.42
2	A	601	SLU	O28-C24	-2.10	1.37	1.43
2	A	601	SLU	C23-C22	-2.05	1.47	1.53

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	SLU	O19-S17-O18	-19.03	101.63	121.45
2	B	601	SLU	O19-S17-O18	-17.13	103.60	121.45
2	A	601	SLU	C8-C11-S12	16.55	129.48	122.35
2	B	601	SLU	C8-C11-S12	13.01	127.95	122.35
2	A	601	SLU	C8-N7-C3	9.76	120.99	103.62
2	B	601	SLU	C14-C13-S12	-9.20	101.24	111.97
2	B	601	SLU	C8-N7-C3	9.13	119.88	103.62
2	B	601	SLU	N31-C30-N29	-8.56	121.55	128.71
2	A	601	SLU	C14-C13-S12	-8.26	102.33	111.97
2	A	601	SLU	N31-C30-N29	-7.56	122.39	128.71
2	B	601	SLU	C11-C8-S9	7.25	125.47	122.35
2	A	601	SLU	C3-C4-S9	-7.21	102.97	112.03
2	B	601	SLU	C13-C14-N15	7.20	120.65	109.74
2	B	601	SLU	C3-C4-S9	-6.69	103.62	112.03
2	A	601	SLU	C13-C14-N15	6.48	119.57	109.74
2	A	601	SLU	C11-C8-S9	6.06	124.96	122.35
2	A	601	SLU	C5-C4-S9	5.26	134.69	125.62
2	B	601	SLU	C5-C4-S9	4.86	134.00	125.62
2	A	601	SLU	C14-N15-C11	4.67	117.14	104.57
2	A	601	SLU	N31-C32-N35	4.62	133.78	125.43
2	B	601	SLU	C14-N15-C11	4.55	116.80	104.57
2	A	601	SLU	S9-C8-N7	-4.48	104.65	117.14
2	B	601	SLU	C22-O26-C25	-4.25	105.13	109.75
2	B	601	SLU	N31-C32-N35	4.10	132.83	125.43
2	B	601	SLU	S9-C8-N7	-4.02	105.93	117.14
2	B	601	SLU	O26-C25-N35	3.65	111.83	108.44
2	B	601	SLU	O20-S17-N40	3.61	113.44	106.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	SLU	C32-C33-N37	-3.58	106.46	109.52
2	A	601	SLU	C22-O26-C25	-3.53	105.92	109.75
2	A	601	SLU	C8-S9-C4	3.19	95.79	90.16
2	B	601	SLU	C4-C3-N7	3.16	115.78	108.37
2	B	601	SLU	O39-C16-N40	3.08	124.85	121.09
2	A	601	SLU	C4-C3-N7	3.09	115.60	108.37
2	A	601	SLU	C11-C8-N7	3.06	130.39	122.38
2	B	601	SLU	C2-C3-C4	-2.99	116.21	121.01
2	B	601	SLU	C13-S12-C11	2.98	96.40	90.79
2	A	601	SLU	O20-S17-N40	2.92	112.11	106.55
2	A	601	SLU	C33-C32-N31	-2.91	119.37	125.70
2	A	601	SLU	C14-C16-N40	2.80	116.18	113.76
2	B	601	SLU	O19-S17-N40	2.78	112.37	108.33
2	B	601	SLU	C32-C33-N37	-2.74	107.17	109.52
2	A	601	SLU	O20-S17-O19	2.73	114.19	105.65
2	A	601	SLU	C2-C3-C4	-2.72	116.64	121.01
2	B	601	SLU	C8-S9-C4	2.62	94.78	90.16
2	A	601	SLU	C13-S12-C11	2.58	95.64	90.79
2	A	601	SLU	O19-S17-N40	2.46	111.91	108.33
2	B	601	SLU	C11-C8-N7	2.37	128.60	122.38
2	A	601	SLU	O26-C22-C23	2.33	109.88	105.17
2	B	601	SLU	C21-O20-S17	2.31	122.66	118.28
2	B	601	SLU	O20-S17-O19	2.30	112.83	105.65
2	A	601	SLU	C30-N31-C32	2.24	120.38	114.01
2	B	601	SLU	O18-S17-N40	-2.23	105.11	108.33
2	B	601	SLU	C33-C32-N31	-2.19	120.93	125.70
2	B	601	SLU	C5-C4-C3	2.19	122.36	120.82
2	A	601	SLU	C5-C4-C3	2.16	122.34	120.82
2	B	601	SLU	C25-N35-C32	-2.13	122.96	126.64
2	B	601	SLU	O26-C25-C24	2.10	110.00	106.77
2	B	601	SLU	C30-N31-C32	2.09	119.96	114.01
2	B	601	SLU	C30-N29-C34	2.08	122.53	118.77
2	B	601	SLU	C16-N40-S17	-2.05	120.84	123.54
2	A	601	SLU	C36-N37-C33	2.00	109.80	103.58

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	SLU	O39-C16-C14-N15
2	B	601	SLU	O39-C16-C14-N15

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 ⓘ

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	522/555 (94%)	-0.15	5 (0%)	79 80	17, 28, 67, 84	0
1	B	543/555 (97%)	-0.22	1 (0%)	93 94	18, 33, 55, 72	0
All	All	1065/1110 (95%)	-0.19	6 (0%)	86 88	17, 31, 63, 84	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	514	GLY	3.5
1	A	520	ASP	3.4
1	A	539	LEU	2.9
1	A	457	ILE	2.5
1	B	175	GLY	2.0
1	A	175	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SLU	A	601	40/40	0.14	-0.07	18,23,27,34	0
2	SLU	B	601	40/40	0.13	-0.45	21,27,30,35	0

6.5 Other polymers ⓘ

There are no such residues in this entry.