



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:46 PM GMT

PDB ID : 4G37
Title : Structure of cross-linked firefly luciferase in second catalytic conformation
Authors : Sundlov, J.A.; Branchini, B.R.; Gulick, A.M.
Deposited on : 2012-07-13
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

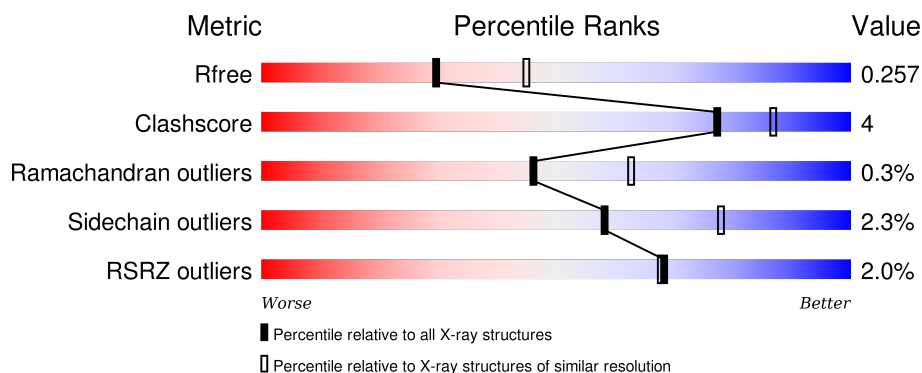
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	555	<div> <div>2%</div> <div>84%</div> <div>11%</div> <div>• •</div> </div>
1	B	555	<div> <div>2%</div> <div>88%</div> <div>9%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	A	603	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8395 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 Luciferin 4-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	533	Total	C	N	O	S	0	0	0
			4015	2583	668	749	15			
1	B	539	Total	C	N	O	S	0	0	0
			4069	2623	674	757	15			

There are 32 discrepancies between the modelled and reference sequences:

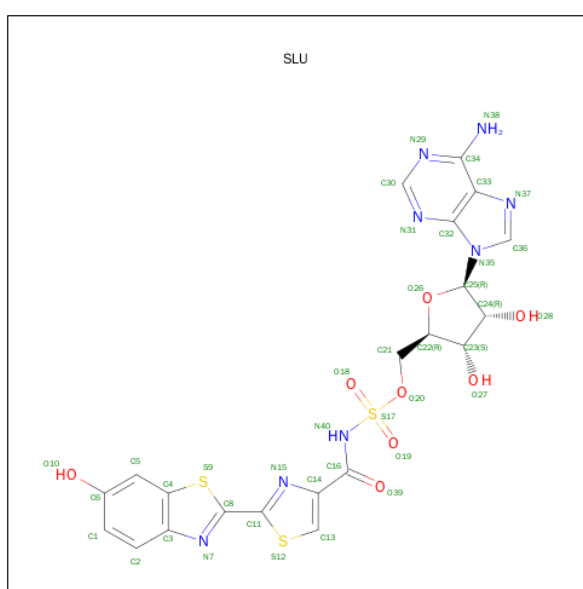
Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP P08659
A	-3	PRO	-	EXPRESSION TAG	UNP P08659
A	-2	LEU	-	EXPRESSION TAG	UNP P08659
A	-1	GLY	-	EXPRESSION TAG	UNP P08659
A	0	SER	-	EXPRESSION TAG	UNP P08659
A	81	SER	CYS	ENGINEERED MUTATION	UNP P08659
A	108	CYS	ILE	ENGINEERED MUTATION	UNP P08659
A	214	ALA	THR	ENGINEERED MUTATION	UNP P08659
A	215	LEU	ALA	ENGINEERED MUTATION	UNP P08659
A	216	ALA	CYS	ENGINEERED MUTATION	UNP P08659
A	232	ALA	ILE	ENGINEERED MUTATION	UNP P08659
A	258	SER	CYS	ENGINEERED MUTATION	UNP P08659
A	295	LEU	PHE	ENGINEERED MUTATION	UNP P08659
A	354	LYS	GLU	ENGINEERED MUTATION	UNP P08659
A	391	SER	CYS	ENGINEERED MUTATION	UNP P08659
A	447	CYS	TYR	ENGINEERED MUTATION	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
B	81	SER	CYS	ENGINEERED MUTATION	UNP P08659
B	108	CYS	ILE	ENGINEERED MUTATION	UNP P08659
B	214	ALA	THR	ENGINEERED MUTATION	UNP P08659
B	215	LEU	ALA	ENGINEERED MUTATION	UNP P08659

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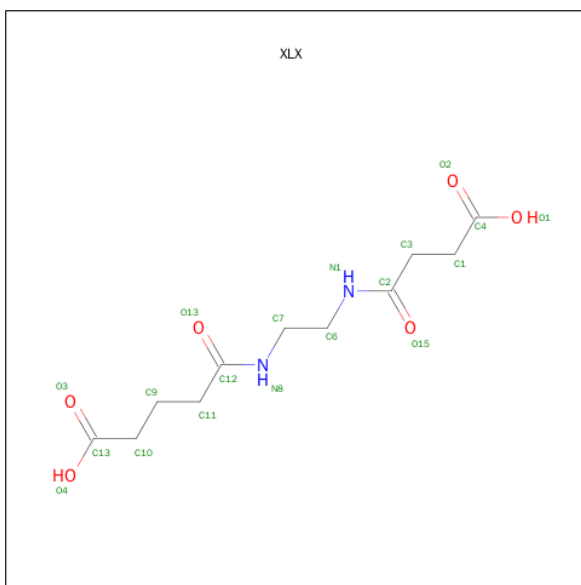
Chain	Residue	Modelled	Actual	Comment	Reference
B	216	ALA	CYS	ENGINEERED MUTATION	UNP P08659
B	232	ALA	ILE	ENGINEERED MUTATION	UNP P08659
B	258	SER	CYS	ENGINEERED MUTATION	UNP P08659
B	295	LEU	PHE	ENGINEERED MUTATION	UNP P08659
B	354	LYS	GLU	ENGINEERED MUTATION	UNP P08659
B	391	SER	CYS	ENGINEERED MUTATION	UNP P08659
B	447	CYS	TYR	ENGINEERED MUTATION	UNP P08659

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



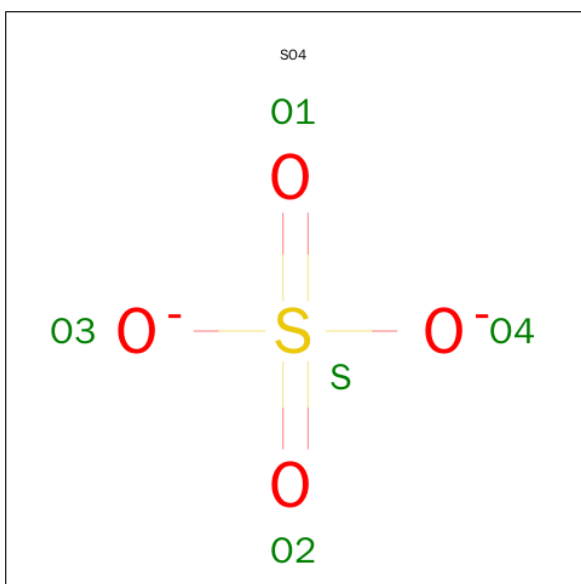
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			40	21	8	8	3		
2	B	1	Total	C	N	O	S	0	0
			40	21	8	8	3		

- Molecule 3 is 4,4'-(ETHYLENEDIIMINO)BIS[4-OXOBUTYRATE] (three-letter code: XLX) (formula: C₁₁H₁₈N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	6	2	2		
3	B	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		

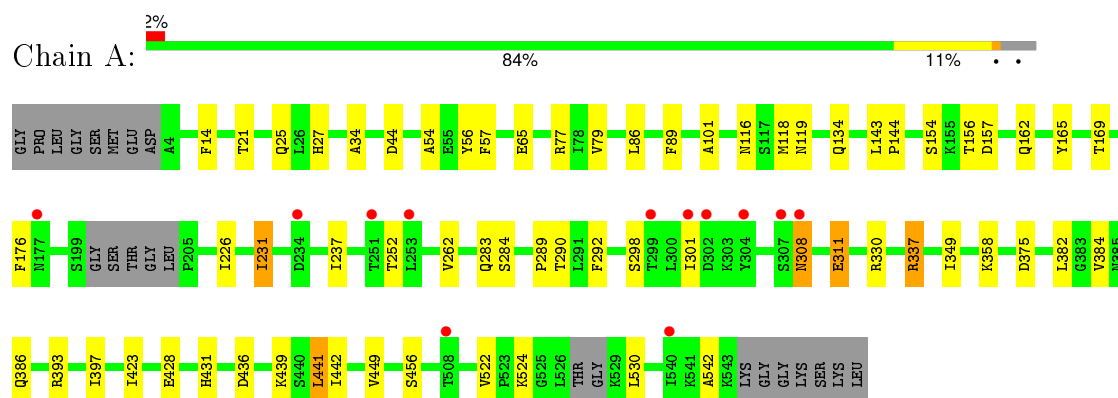
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	78	Total	O	0	0
			78	78		
6	B	112	Total	O	0	0
			112	112		

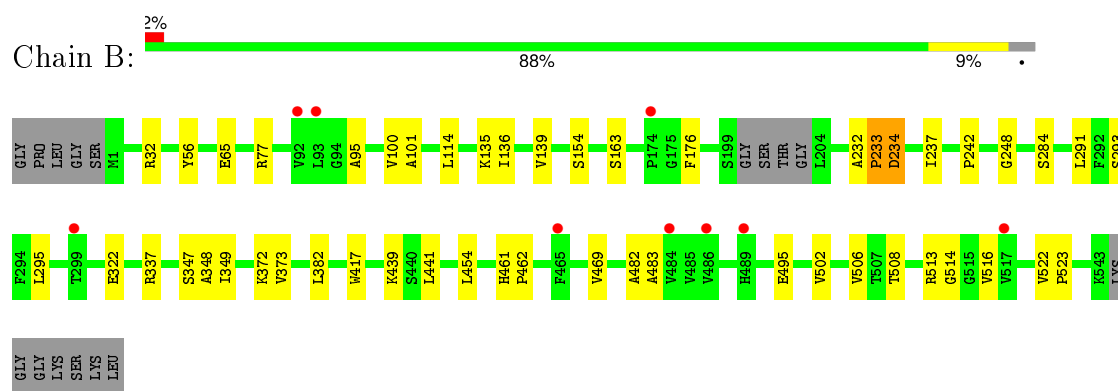
3 Residue-property plots [i](#)

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.

• Molecule 1: Luciferin 4-monooxygenase



• Molecule 1: Luciferin 4-monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	75.64Å 184.09Å 170.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.22 – 2.40 29.22 – 2.40	Depositor EDS
% Data completeness (in resolution range)	94.1 (29.22-2.40) 94.1 (29.22-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_837)	Depositor
R, R_{free}	0.199 , 0.259 0.198 , 0.257	Depositor DCC
R_{free} test set	2275 reflections (5.40%)	DCC
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 31.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 44401 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8395	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SLU, CL, SO4, XLX

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	0.40	0/4102	0.55	0/5581
1	B	0.42	0/4159	0.57	0/5659
All	All	0.41	0/8261	0.56	0/11240

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

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	4015	0	3906	36	0
1	B	4069	0	3980	25	0
2	A	40	0	18	0	0
2	B	40	0	18	0	0
3	A	10	0	6	0	0
3	B	10	0	6	0	0
4	A	10	0	0	0	0
4	B	10	0	0	0	0
5	B	1	0	0	0	0
6	A	78	0	0	0	0
6	B	112	0	0	0	0
All	All	8395	0	7934	60	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 4.

All (60) 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:A:237:ILE:HG22	1:A:284:SER:HB3	1.72	0.71
1:B:77:ARG:HG2	1:B:101:ALA:HB3	1.72	0.70
1:A:231:ILE:H	1:A:231:ILE:HD12	1.59	0.67
1:B:232:ALA:O	1:B:234:ASP:N	2.29	0.64
1:B:373:VAL:HG23	1:B:382:LEU:HB2	1.81	0.62
1:B:135:LYS:O	1:B:139:VAL:HG23	2.00	0.62
1:B:114:LEU:HD21	1:B:136:ILE:HD13	1.82	0.61
1:A:393:ARG:HA	1:A:397:ILE:HD12	1.84	0.60
1:B:495:GLU:HG3	1:B:516:VAL:HB	1.84	0.60
1:A:439:LYS:HE3	1:A:441:LEU:HD22	1.85	0.59
1:B:114:LEU:HD23	1:B:139:VAL:HG21	1.86	0.57
1:A:375:ASP:HB2	1:A:382:LEU:HD21	1.88	0.56
1:A:524:LYS:HG2	1:A:530:LEU:HD23	1.89	0.54
1:A:165:TYR:O	1:A:169:THR:HG23	2.06	0.54
1:A:14:PHE:CD1	1:B:32:ARG:HD3	2.44	0.53
1:B:242:PRO:O	1:B:248:GLY:HA3	2.10	0.52
1:A:77:ARG:HG3	1:A:101:ALA:HB3	1.93	0.51
1:B:482:ALA:HB2	1:B:514:GLY:HA3	1.94	0.50
1:B:502:VAL:O	1:B:506:VAL:HG22	2.11	0.50
1:B:508:THR:O	1:B:513:ARG:NH1	2.40	0.50
1:A:522:VAL:O	1:A:524:LYS:HG3	2.13	0.48
1:B:237:ILE:HG22	1:B:284:SER:HB3	1.96	0.48
1:A:65:GLU:HG2	1:A:176:PHE:CE1	2.48	0.48
1:B:439:LYS:HE3	1:B:441:LEU:HD22	1.96	0.47
1:B:522:VAL:HA	1:B:523:PRO:HD3	1.80	0.47
1:A:27:HIS:CD2	1:A:57:PHE:HE1	2.33	0.47
1:A:298:SER:O	1:A:330:ARG:NH1	2.48	0.46
1:A:89:PHE:CE1	1:A:252:THR:HG21	2.51	0.46
1:B:291:LEU:O	1:B:295:LEU:HB2	2.15	0.46
1:A:226:ILE:O	1:A:337:ARG:NH2	2.49	0.46
1:B:469:VAL:HG22	1:B:483:ALA:HB2	1.97	0.46
1:A:21:THR:O	1:A:25:GLN:HG3	2.16	0.45
1:A:143:LEU:HA	1:A:143:LEU:HD12	1.70	0.45
1:B:372:LYS:HE2	1:B:417:TRP:CD2	2.52	0.45
1:A:311:GLU:OE1	1:A:337:ARG:HD3	2.17	0.44
1:A:237:ILE:HD11	1:A:262:VAL:HG13	1.99	0.44
1:B:65:GLU:HG2	1:B:176:PHE:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:ILE:HD12	1:A:436:ASP:HB3	1.99	0.44
1:B:295:LEU:HA	1:B:295:LEU:HD12	1.84	0.44
1:A:384:VAL:HG21	1:A:428:GLU:HG3	2.00	0.43
1:A:442:ILE:HB	1:A:449:VAL:HB	2.00	0.43
1:A:79:VAL:HG21	1:A:118:MET:HE3	2.01	0.43
1:A:118:MET:HB3	1:A:118:MET:HE2	1.82	0.43
1:B:154:SER:O	1:B:163:SER:HB2	2.18	0.43
1:A:289:PRO:HA	1:A:292:PHE:CD1	2.53	0.43
1:A:157:ASP:OD1	1:A:162:GLN:HA	2.19	0.43
1:A:34:ALA:HB1	1:A:54:ALA:HA	2.01	0.43
1:A:393:ARG:HA	1:A:397:ILE:CD1	2.48	0.42
1:A:298:SER:HB3	1:A:301:ILE:HD11	2.01	0.42
1:A:143:LEU:HA	1:A:144:PRO:HD2	1.87	0.42
1:A:154:SER:OG	1:A:156:THR:O	2.34	0.42
1:A:44:ASP:N	1:A:86:LEU:HD21	2.35	0.42
1:B:454:LEU:HD12	1:B:469:VAL:HG11	2.03	0.41
1:A:358:LYS:HD2	1:A:431:HIS:ND1	2.35	0.41
1:A:283:GLN:HA	1:A:308:ASN:O	2.20	0.41
1:A:284:SER:OG	1:A:311:GLU:HG2	2.20	0.41
1:A:27:HIS:CD2	1:A:57:PHE:CE1	3.08	0.41
1:B:347:SER:OG	1:B:348:ALA:N	2.52	0.40
1:B:95:ALA:HB1	1:B:100:VAL:O	2.22	0.40
1:B:461:HIS:HA	1:B:462:PRO:HD3	1.91	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	527/555 (95%)	508 (96%)	18 (3%)	1 (0%)	52 69
1	B	535/555 (96%)	515 (96%)	18 (3%)	2 (0%)	39 56

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1062/1110 (96%)	1023 (96%)	36 (3%)	3 (0%)	46 63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	233	PRO
1	A	542	ALA
1	B	234	ASP

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	412/464 (89%)	399 (97%)	13 (3%)	46 68
1	B	419/464 (90%)	413 (99%)	6 (1%)	74 88
All	All	831/928 (90%)	812 (98%)	19 (2%)	58 78

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	TYR
1	A	116	ASN
1	A	119	ASN
1	A	134	GLN
1	A	231	ILE
1	A	290	THR
1	A	308	ASN
1	A	311	GLU
1	A	337	ARG
1	A	349	ILE
1	A	386	GLN
1	A	441	LEU
1	A	456	SER
1	B	56	TYR
1	B	233	PRO

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Mol	Chain	Res	Type
1	B	293	SER
1	B	322	GLU
1	B	337	ARG
1	B	349	ILE

Some 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](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	SLU	A	601	-	35,45,45	1.83	9 (25%)	37,68,68	3.23	7 (18%)
3	XLX	A	602	1	9,9,18	0.27	0	10,10,21	0.60	0
4	SO4	A	603	-	4,4,4	0.09	0	6,6,6	0.14	0
4	SO4	A	604	-	4,4,4	0.15	0	6,6,6	0.32	0
2	SLU	B	601	-	35,45,45	1.68	10 (28%)	37,68,68	3.64	12 (32%)
3	XLX	B	602	1	9,9,18	0.24	0	10,10,21	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	B	604	-	4,4,4	0.18	0	6,6,6	0.11	0
4	SO4	B	605	-	4,4,4	0.18	0	6,6,6	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	SLU	A	601	-	-	0/10/39/39	0/6/6/6
3	XLX	A	602	1	-	4/7/7/18	0/0/0/0
4	SO4	A	603	-	-	0/0/0/0	0/0/0/0
4	SO4	A	604	-	-	0/0/0/0	0/0/0/0
2	SLU	B	601	-	-	0/10/39/39	0/6/6/6
3	XLX	B	602	1	-	4/7/7/18	0/0/0/0
4	SO4	B	604	-	-	0/0/0/0	0/0/0/0
4	SO4	B	605	-	-	0/0/0/0	0/0/0/0

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	SLU	C24-C23	-2.51	1.46	1.53
2	B	601	SLU	C24-C23	-2.44	1.46	1.53
2	A	601	SLU	C14-N15	-2.42	1.30	1.37
2	B	601	SLU	O28-C24	-2.42	1.37	1.43
2	A	601	SLU	C3-N7	-2.15	1.31	1.39
2	B	601	SLU	O26-C22	-2.14	1.40	1.45
2	B	601	SLU	C14-N15	-2.12	1.31	1.37
2	B	601	SLU	C3-N7	-2.10	1.32	1.39
2	A	601	SLU	O28-C24	-2.03	1.38	1.43
2	B	601	SLU	C5-C4	2.02	1.40	1.37
2	A	601	SLU	C5-C4	2.15	1.41	1.37
2	B	601	SLU	C13-S12	2.62	1.74	1.70
2	A	601	SLU	C13-S12	2.73	1.74	1.70
2	A	601	SLU	C34-N38	3.09	1.44	1.34
2	B	601	SLU	C34-N38	3.13	1.44	1.34
2	B	601	SLU	O19-S17	3.17	1.45	1.42
2	A	601	SLU	O19-S17	3.83	1.45	1.42
2	B	601	SLU	C16-N40	4.10	1.43	1.39
2	A	601	SLU	C16-N40	5.53	1.45	1.39

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	SLU	O19-S17-O18	-12.66	105.61	120.77
2	A	601	SLU	O19-S17-O18	-11.14	107.42	120.77
2	B	601	SLU	N31-C30-N29	-9.24	121.82	128.89
2	A	601	SLU	N31-C30-N29	-7.92	122.83	128.89
2	A	601	SLU	C14-C13-S12	-7.40	102.70	111.79
2	B	601	SLU	C14-C13-S12	-7.32	102.80	111.79
2	B	601	SLU	C16-N40-S17	-6.89	115.38	123.35
2	A	601	SLU	C22-O26-C25	-4.94	104.29	109.72
2	B	601	SLU	C22-O26-C25	-4.73	104.52	109.72
2	B	601	SLU	C32-C33-N37	-3.17	106.56	109.48
2	A	601	SLU	C32-C33-N37	-2.92	106.79	109.48
2	B	601	SLU	C24-C25-N35	-2.74	110.10	114.29
2	B	601	SLU	C13-C14-C16	-2.08	115.10	126.35
2	B	601	SLU	O39-C16-C14	2.15	125.80	121.23
2	B	601	SLU	O20-S17-O19	2.27	112.69	105.66
2	B	601	SLU	C4-C3-N7	3.21	115.60	108.16
2	A	601	SLU	C4-C3-N7	3.24	115.68	108.16
2	B	601	SLU	C8-N7-C3	7.58	118.70	103.83
2	A	601	SLU	C8-N7-C3	8.32	120.15	103.83

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	XLX	C11-C12-N8-C7
3	A	602	XLX	C3-C2-N1-C6
3	B	602	XLX	C11-C12-N8-C7
3	B	602	XLX	C3-C2-N1-C6
3	B	602	XLX	O15-C2-N1-C6
3	A	602	XLX	O15-C2-N1-C6
3	B	602	XLX	O13-C12-N8-C7
3	A	602	XLX	O13-C12-N8-C7

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	533/555 (96%)	-0.03	12 (2%) 64 63	21, 42, 71, 95	0
1	B	539/555 (97%)	-0.13	9 (1%) 73 72	21, 38, 65, 88	0
All	All	1072/1110 (96%)	-0.08	21 (1%) 68 68	21, 40, 69, 95	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	486	VAL	4.2
1	A	302	ASP	4.0
1	A	301	ILE	3.4
1	B	299	THR	3.2
1	B	484	VAL	3.2
1	B	92	VAL	3.2
1	A	234	ASP	3.1
1	B	489	HIS	2.6
1	B	465	PHE	2.5
1	A	304	TYR	2.5
1	B	174	PRO	2.5
1	B	517	VAL	2.4
1	A	253	LEU	2.3
1	A	308	ASN	2.3
1	B	93	LEU	2.2
1	A	307	SER	2.1
1	A	251	THR	2.1
1	A	508	THR	2.1
1	A	540	ILE	2.1
1	A	177	ASN	2.0
1	A	299	THR	2.0

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

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

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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	SO4	A	603	5/5	0.92	0.18	4.48	54,58,70,73	0
3	XLX	B	602	10/19	0.85	0.17	1.35	60,68,77,77	0
3	XLX	A	602	10/19	0.84	0.20	0.66	68,70,77,78	0
2	SLU	A	601	40/40	0.97	0.14	-0.18	22,28,31,33	0
2	SLU	B	601	40/40	0.98	0.12	-0.61	22,27,31,33	0
4	SO4	A	604	5/5	0.96	0.14	-	49,61,65,70	0
4	SO4	B	605	5/5	0.96	0.18	-	37,46,69,71	0
5	CL	B	603	1/1	0.82	0.06	-	76,76,76,76	0
4	SO4	B	604	5/5	0.97	0.08	-	46,46,51,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.