



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:52 PM GMT

PDB ID : 4MX2
Title : Crystal Structure of adenylosuccinate lyase from Leishmania donovani
Authors : Wernimont, A.K.; Loppnau, P.; Dong, A.; Krojer, T.; Bradley, A.; Bushell, S.; von Delft, F.; Robinson, D.; Gilbert, I.; Bountra, C.; Arrowsmith, C.H.; Edwards, A.M.; Hui, R.; Mottaghi, K.; Structural Genomics Consortium (SGC)
Deposited on : 2013-09-25
Resolution : 1.90 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

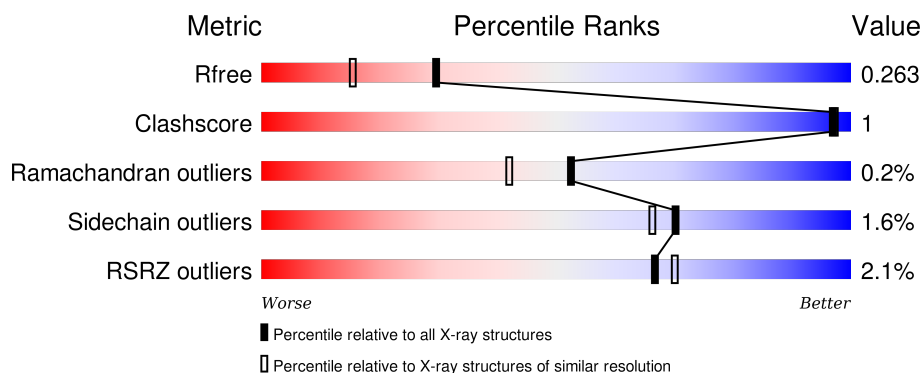
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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	480	<div> <div>2%</div> <div> <div></div> <div>85%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	480	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>5%</div> <div>9%</div> </div> </div>
1	C	480	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>•</div> <div>11%</div> </div> </div>
1	D	480	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>•</div> <div>9%</div> </div> </div>
1	E	480	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>•</div> <div>11%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	480	 3% 85% 5% 10%
1	G	480	 3% 85% 11%
1	H	480	 2% 86% 5% 9%

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	GOL	E	501	-	-	X	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 29460 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 Adenylosuccinate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	8	0
			3333	2145	557	617	14			
1	B	439	Total	C	N	O	S	0	15	0
			3522	2261	595	652	14			
1	C	426	Total	C	N	O	S	0	8	0
			3367	2167	568	618	14			
1	D	439	Total	C	N	O	S	0	9	0
			3487	2239	588	647	13			
1	E	425	Total	C	N	O	S	0	9	0
			3347	2154	560	619	14			
1	F	431	Total	C	N	O	S	0	12	0
			3450	2222	577	637	14			
1	G	427	Total	C	N	O	S	0	9	0
			3375	2174	567	620	14			
1	H	438	Total	C	N	O	S	0	14	0
			3490	2247	584	644	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
B	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
C	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
D	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
E	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
F	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
G	0	GLY	-	EXPRESSION TAG	UNP A7LBL3
H	0	GLY	-	EXPRESSION TAG	UNP A7LBL3

- Molecule 2 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	E	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	F	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	G	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	H	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	C	O	0	0
			6	3	3		

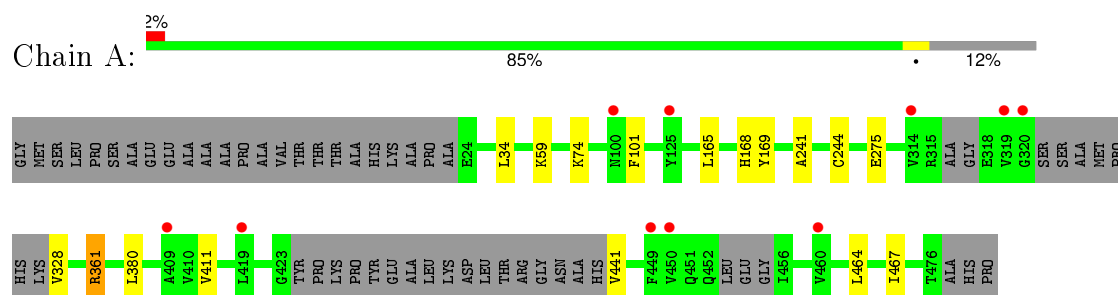
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	232	Total 233	O 233	0	1
5	B	244	Total 244	O 244	0	0
5	C	212	Total 212	O 212	0	0
5	D	264	Total 265	O 265	0	1
5	E	215	Total 215	O 215	0	0
5	F	239	Total 239	O 239	0	0
5	G	217	Total 217	O 217	0	0
5	H	259	Total 259	O 259	0	0

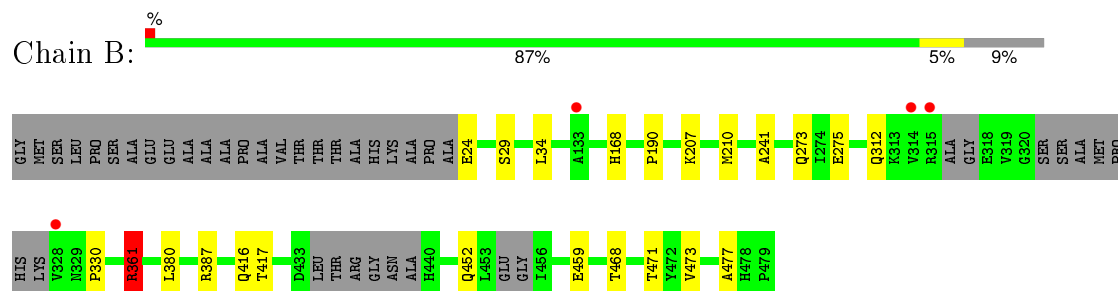
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

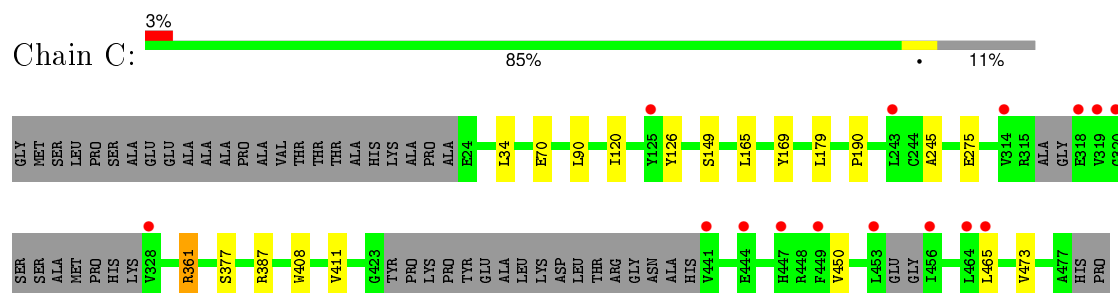
- Molecule 1: Adenylosuccinate lyase



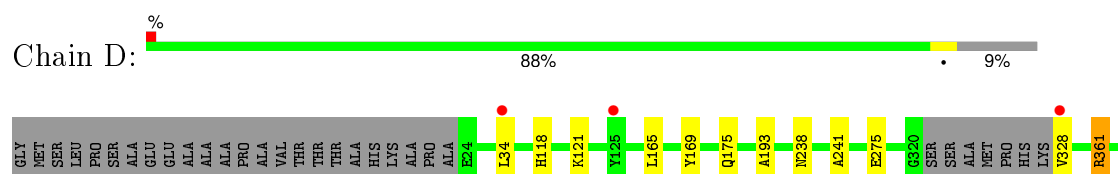
- Molecule 1: Adenylosuccinate lyase



- Molecule 1: Adenylosuccinate lyase



- Molecule 1: Adenylosuccinate lyase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	124.02Å 194.70Å 147.51Å 90.00° 91.31° 90.00°	Depositor
Resolution (Å)	34.98 – 1.90 34.86 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.7 (34.98-1.90) 89.4 (34.86-1.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 1.89Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.218 , 0.253 0.226 , 0.263	Depositor DCC
R_{free} test set	13161 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	16.5	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.9	EDS
Estimated twinning fraction	0.199 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	6 of 261938 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	29460	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.58 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.4942e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, AMP

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	0.50	0/3424	0.59	0/4649
1	B	0.51	0/3626	0.59	0/4921
1	C	0.50	0/3460	0.59	0/4692
1	D	0.50	0/3585	0.59	0/4870
1	E	0.48	0/3442	0.59	0/4673
1	F	0.50	0/3556	0.59	0/4827
1	G	0.49	0/3471	0.59	0/4710
1	H	0.50	0/3607	0.59	0/4900
All	All	0.50	0/28171	0.59	0/38242

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	3333	0	3288	7	0
1	B	3522	0	3471	9	0
1	C	3367	0	3322	9	0
1	D	3487	0	3454	10	0
1	E	3347	0	3299	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3450	0	3429	13	0
1	G	3375	0	3335	6	0
1	H	3490	0	3460	13	0
2	A	23	0	12	0	0
2	B	23	0	12	0	0
2	C	23	0	12	0	0
2	D	23	0	12	0	0
2	E	23	0	12	0	0
2	F	23	0	12	0	0
2	G	23	0	12	0	0
2	H	23	0	12	0	0
3	C	5	0	0	0	0
3	G	5	0	0	0	0
3	H	5	0	0	0	0
4	E	6	0	8	4	0
5	A	233	0	0	0	0
5	B	244	0	0	0	0
5	C	212	0	0	0	0
5	D	265	0	0	1	0
5	E	215	0	0	0	0
5	F	239	0	0	2	0
5	G	217	0	0	0	0
5	H	259	0	0	0	0
All	All	29460	0	27162	70	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 1.

All (70) 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:E:157:ILE:HD11	4:E:501:GOL:H32	1.39	1.05
1:A:244[B]:CYS:HG	1:E:472:TYR:HD2	1.05	0.98
1:H:170:ILE:HD11	1:H:223:LEU:HD23	1.67	0.76
1:E:366:SER:HA	4:E:501:GOL:H12	1.67	0.75
1:E:157:ILE:CD1	4:E:501:GOL:H32	2.18	0.72
1:B:417[B]:THR:HG21	1:F:241:ALA:HB1	1.72	0.71
1:F:259[B]:LEU:HD22	1:F:265:LEU:HB2	1.77	0.66
1:B:241:ALA:HB1	1:F:417[B]:THR:HG21	1.77	0.65
1:F:259[B]:LEU:HD23	1:F:263:LEU:HD12	1.78	0.65
1:D:241:ALA:HB1	1:H:417[B]:THR:HG21	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:155:THR:HG22	1:F:259[B]:LEU:HD21	1.81	0.62
1:E:118:HIS:HD2	1:E:121:LYS:H	1.48	0.60
1:B:468:THR:H	1:B:471[A]:THR:HG22	1.65	0.60
1:A:59:LYS:HD3	1:A:101:PHE:HB3	1.87	0.57
1:A:241:ALA:HB1	1:E:417[B]:THR:HG21	1.88	0.56
1:H:170:ILE:HD11	1:H:223:LEU:CD2	2.37	0.54
1:G:190:PRO:HB3	1:G:473:VAL:HG11	1.89	0.54
1:D:415:ILE:HG23	1:D:464:LEU:HD22	1.89	0.54
1:H:120:ILE:HG13	1:H:149:SER:HB3	1.90	0.54
1:D:118:HIS:HD2	1:D:121:LYS:H	1.55	0.52
1:E:165:LEU:HA	1:E:169:TYR:HB3	1.92	0.50
1:H:118[A]:HIS:HD2	1:H:121:LYS:H	1.59	0.49
1:F:259[B]:LEU:CD2	1:F:263:LEU:HD12	2.43	0.49
1:E:105:ASP:HA	1:E:108:ILE:HD12	1.95	0.49
1:E:60:TYR:HB3	1:E:156:SER:HB2	1.94	0.49
1:B:190:PRO:HB3	1:B:473:VAL:HG11	1.94	0.49
1:F:342:MET:HA	1:F:342:MET:HE3	1.94	0.48
1:B:210[B]:MET:SD	1:B:477:ALA:HA	2.54	0.48
1:F:190:PRO:HB3	1:F:473:VAL:HG11	1.96	0.48
1:D:165:LEU:HA	1:D:169:TYR:HB3	1.96	0.47
1:C:377:SER:OG	1:D:34:LEU:HD11	2.13	0.47
1:B:312:GLN:HE22	1:B:330:PRO:HA	1.79	0.47
1:H:312:GLN:HE22	1:H:330:PRO:HA	1.80	0.47
1:H:210[B]:MET:SD	1:H:477:ALA:HA	2.54	0.47
1:H:190:PRO:HB3	1:H:473:VAL:HG11	1.97	0.47
1:B:273:GLN:HB3	1:B:361:ARG:HG2	1.97	0.47
1:C:245:ALA:HB2	1:G:417:THR:HG23	1.98	0.46
1:H:175:GLN:HB3	1:H:387:ARG:NH2	2.31	0.45
1:A:168[B]:HIS:HD2	1:A:380:LEU:HD11	1.81	0.45
1:C:165:LEU:HA	1:C:169:TYR:HB3	1.98	0.45
1:E:366:SER:HA	4:E:501:GOL:C1	2.42	0.44
1:A:411:VAL:HB	1:A:467:ILE:HG22	1.99	0.44
1:C:450:VAL:HG11	1:C:465:LEU:HG	1.99	0.44
1:F:252:LEU:HD21	5:F:605:HOH:O	2.18	0.44
1:H:419:LEU:HB3	1:H:424:TYR:HB3	2.00	0.44
1:D:241:ALA:O	1:H:417[A]:THR:HG21	2.17	0.44
1:G:105:ASP:HA	1:G:108:ILE:HD12	2.00	0.43
1:D:118:HIS:CD2	1:D:121:LYS:H	2.36	0.43
1:G:165:LEU:HA	1:G:169:TYR:HB3	2.00	0.43
1:A:165:LEU:HA	1:A:169:TYR:HB3	2.00	0.43
1:E:118:HIS:CD2	1:E:121:LYS:H	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:408:TRP:O	1:H:411:VAL:HG22	2.18	0.43
1:E:157:ILE:HA	1:E:160:LEU:HG	2.00	0.42
1:C:179:LEU:HD22	1:C:387[B]:ARG:HG3	2.02	0.42
1:F:60:TYR:HB3	1:F:156[B]:SER:HB2	2.00	0.42
1:C:120:ILE:HG21	1:C:149:SER:HA	2.02	0.42
1:B:207:LYS:NZ	5:F:766:HOH:O	2.53	0.42
1:G:63:GLN:HA	1:G:98:PHE:CD2	2.54	0.42
1:C:408:TRP:O	1:C:411:VAL:HG22	2.20	0.41
1:G:450:VAL:HG11	1:G:465:LEU:HG	2.02	0.41
1:A:464:LEU:HA	1:A:467:ILE:HD12	2.03	0.41
1:D:238:ASN:ND2	5:D:852:HOH:O	2.53	0.41
1:B:168[B]:HIS:HD2	1:B:380:LEU:HD11	1.85	0.41
1:F:137:GLU:O	1:F:140:LYS:HG2	2.20	0.41
1:C:190:PRO:HB3	1:C:473:VAL:HG11	2.02	0.40
1:F:159:MET:HG3	1:F:265:LEU:HD11	2.03	0.40
1:C:70:GLU:HG3	1:C:90:LEU:HD13	2.04	0.40
1:F:118:HIS:CD2	1:F:121:LYS:H	2.39	0.40
1:D:193:ALA:CB	1:H:236:THR:HG21	2.51	0.40
1:D:408:TRP:O	1:D:411:VAL:HG22	2.20	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	422/480 (88%)	409 (97%)	12 (3%)	1 (0%)	52 42
1	B	444/480 (92%)	422 (95%)	21 (5%)	1 (0%)	52 42
1	C	424/480 (88%)	408 (96%)	15 (4%)	1 (0%)	52 42
1	D	441/480 (92%)	421 (96%)	19 (4%)	1 (0%)	52 42
1	E	424/480 (88%)	408 (96%)	15 (4%)	1 (0%)	52 42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	433/480 (90%)	415 (96%)	17 (4%)	1 (0%)	52	42
1	G	427/480 (89%)	411 (96%)	15 (4%)	1 (0%)	52	42
1	H	442/480 (92%)	422 (96%)	19 (4%)	1 (0%)	52	42
All	All	3457/3840 (90%)	3316 (96%)	133 (4%)	8 (0%)	52	42

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	361	ARG
1	B	361	ARG
1	C	361	ARG
1	D	361	ARG
1	E	361	ARG
1	F	361	ARG
1	G	361	ARG
1	H	361	ARG

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	350/403 (87%)	344 (98%)	6 (2%)	68	64
1	B	371/403 (92%)	361 (97%)	10 (3%)	52	43
1	C	350/403 (87%)	346 (99%)	4 (1%)	80	79
1	D	370/403 (92%)	365 (99%)	5 (1%)	74	71
1	E	351/403 (87%)	346 (99%)	5 (1%)	74	71
1	F	368/403 (91%)	362 (98%)	6 (2%)	70	66
1	G	353/403 (88%)	347 (98%)	6 (2%)	68	64
1	H	372/403 (92%)	367 (99%)	5 (1%)	76	73
All	All	2885/3224 (90%)	2838 (98%)	47 (2%)	70	66

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	A	74	LYS
1	A	275	GLU
1	A	328	VAL
1	A	361	ARG
1	A	441	VAL
1	B	24	GLU
1	B	29	SER
1	B	34	LEU
1	B	275	GLU
1	B	361	ARG
1	B	387[A]	ARG
1	B	387[B]	ARG
1	B	416	GLN
1	B	452	GLN
1	B	459	GLU
1	C	34	LEU
1	C	126	TYR
1	C	275	GLU
1	C	361	ARG
1	D	175	GLN
1	D	275	GLU
1	D	328	VAL
1	D	361	ARG
1	D	431	LEU
1	E	34	LEU
1	E	80	THR
1	E	150	GLN
1	E	275	GLU
1	E	361	ARG
1	F	34	LEU
1	F	126	TYR
1	F	275	GLU
1	F	361	ARG
1	F	451	GLN
1	F	459	GLU
1	G	34	LEU
1	G	126	TYR
1	G	275	GLU
1	G	328	VAL
1	G	361	ARG
1	G	422	GLU
1	H	34	LEU

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Mol	Chain	Res	Type
1	H	275	GLU
1	H	328	VAL
1	H	361	ARG
1	H	431	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	167	HIS
1	B	312	GLN
1	B	452	GLN
1	D	166	HIS
1	D	167	HIS
1	D	312	GLN
1	F	452	GLN
1	H	312	GLN

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](#)

12 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	AMP	A	501	-	20,25,25	0.62	0	22,38,38	0.67	0
2	AMP	B	501	-	20,25,25	0.65	0	22,38,38	0.66	0
2	AMP	C	501	-	20,25,25	0.60	0	22,38,38	0.63	0
3	SO4	C	502	-	4,4,4	0.37	0	6,6,6	0.12	0
2	AMP	D	501	-	20,25,25	0.52	0	22,38,38	0.68	0
4	GOL	E	501	-	5,5,5	0.15	0	5,5,5	0.55	0
2	AMP	E	502	-	20,25,25	0.56	0	22,38,38	0.61	0
2	AMP	F	501	-	20,25,25	0.55	0	22,38,38	0.54	0
2	AMP	G	501	-	20,25,25	0.61	0	22,38,38	0.55	0
3	SO4	G	502	-	4,4,4	0.12	0	6,6,6	0.08	0
2	AMP	H	501	-	20,25,25	0.54	0	22,38,38	0.60	0
3	SO4	H	502	-	4,4,4	0.21	0	6,6,6	0.24	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	AMP	A	501	-	-	0/6/26/26	0/3/3/3
2	AMP	B	501	-	-	0/6/26/26	0/3/3/3
2	AMP	C	501	-	-	0/6/26/26	0/3/3/3
3	SO4	C	502	-	-	0/0/0/0	0/0/0/0
2	AMP	D	501	-	-	0/6/26/26	0/3/3/3
4	GOL	E	501	-	-	0/4/4/4	0/0/0/0
2	AMP	E	502	-	-	0/6/26/26	0/3/3/3
2	AMP	F	501	-	-	0/6/26/26	0/3/3/3
2	AMP	G	501	-	-	0/6/26/26	0/3/3/3
3	SO4	G	502	-	-	0/0/0/0	0/0/0/0
2	AMP	H	501	-	-	0/6/26/26	0/3/3/3
3	SO4	H	502	-	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	501	GOL	4	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	424/480 (88%)	0.08	10 (2%) 62 66	9, 17, 43, 83	0
1	B	439/480 (91%)	0.01	4 (0%) 85 87	8, 16, 40, 75	0
1	C	426/480 (88%)	0.17	15 (3%) 48 51	9, 18, 47, 75	0
1	D	439/480 (91%)	0.02	4 (0%) 85 87	8, 16, 38, 64	0
1	E	425/480 (88%)	0.18	13 (3%) 52 56	10, 19, 46, 74	0
1	F	431/480 (89%)	0.08	6 (1%) 78 80	9, 17, 38, 81	0
1	G	427/480 (88%)	0.17	12 (2%) 56 60	9, 19, 43, 77	0
1	H	438/480 (91%)	0.05	9 (2%) 67 70	8, 16, 39, 73	0
All	All	3449/3840 (89%)	0.09	73 (2%) 67 70	8, 17, 43, 83	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	314	VAL	6.9
1	A	314	VAL	6.8
1	B	314	VAL	6.3
1	A	319	VAL	5.8
1	G	314	VAL	5.1
1	H	314	VAL	5.0
1	C	453	LEU	4.3
1	A	419	LEU	4.2
1	A	450	VAL	3.9
1	E	319	VAL	3.9
1	E	449	PHE	3.8
1	E	450	VAL	3.8
1	C	319	VAL	3.7
1	F	328	VAL	3.7
1	A	449	PHE	3.5
1	E	460	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	G	102	ALA	3.2
1	C	449	PHE	3.2
1	H	34	LEU	3.1
1	C	320	GLY	3.1
1	D	34	LEU	3.0
1	F	318	GLU	3.0
1	A	100	ASN	3.0
1	E	419	LEU	3.0
1	C	464	LEU	3.0
1	E	447	HIS	2.8
1	G	419	LEU	2.8
1	F	315	ARG	2.7
1	E	314	VAL	2.7
1	E	100	ASN	2.7
1	E	456	ILE	2.6
1	G	318	GLU	2.6
1	H	315	ARG	2.6
1	E	444	GLU	2.6
1	C	314	VAL	2.6
1	C	444	GLU	2.5
1	A	125	TYR	2.5
1	A	460	VAL	2.4
1	D	328	VAL	2.4
1	D	453	LEU	2.4
1	B	133	ALA	2.4
1	H	430	ALA	2.4
1	H	429	GLU	2.4
1	G	446	VAL	2.3
1	G	69	PHE	2.3
1	H	134[A]	CYS	2.3
1	G	319	VAL	2.3
1	G	441	VAL	2.3
1	C	456	ILE	2.3
1	E	34	LEU	2.3
1	C	465	LEU	2.2
1	H	319	VAL	2.2
1	C	318	GLU	2.2
1	A	409	ALA	2.2
1	D	125	TYR	2.2
1	C	243	LEU	2.2
1	B	328	VAL	2.2
1	F	319	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	456	ILE	2.2
1	H	125	TYR	2.2
1	F	102	ALA	2.2
1	G	447	HIS	2.2
1	C	447	HIS	2.1
1	B	315	ARG	2.1
1	E	446	VAL	2.1
1	E	125	TYR	2.1
1	C	328	VAL	2.1
1	G	100	ASN	2.1
1	A	320	GLY	2.1
1	G	125	TYR	2.0
1	C	441	VAL	2.0
1	G	460	VAL	2.0
1	C	125	TYR	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	GOL	E	501	6/6	0.83	0.17	4.02	26,27,30,31	0
2	AMP	E	502	23/23	0.93	0.12	0.12	15,22,23,24	0
2	AMP	A	501	23/23	0.96	0.09	-0.51	8,14,16,17	0
2	AMP	B	501	23/23	0.97	0.08	-0.55	8,16,19,19	0
2	AMP	D	501	23/23	0.97	0.09	-0.65	10,17,19,20	0
2	AMP	F	501	23/23	0.95	0.09	-0.76	14,19,22,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	AMP	C	501	23/23	0.94	0.09	-0.79	8,14,18,20	0
3	SO4	H	502	5/5	0.94	0.10	-0.82	44,47,48,48	0
3	SO4	C	502	5/5	0.97	0.10	-1.25	29,30,32,32	0
3	SO4	G	502	5/5	0.97	0.07	-1.51	30,34,36,38	0
2	AMP	H	501	23/23	0.96	0.08	-1.52	11,18,20,21	0
2	AMP	G	501	23/23	0.94	0.08	-1.60	11,15,19,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.