



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:00 PM GMT

PDB ID : 1DKA
Title : DIALKYLGLYCINE DECARBOXYLASE STRUCTURE: BIFUNCTIONAL
ACTIVE SITE AND ALKALI METAL BINDING SITES
Authors : Toney, M.D.; Hohenester, E.; Jansonius, J.N.
Deposited on : 1993-06-18
Resolution : 2.60 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

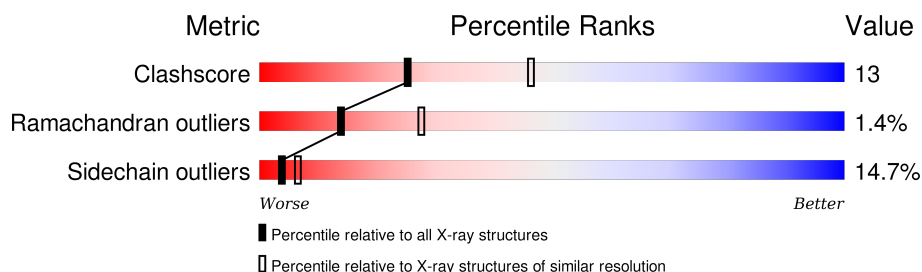
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	433	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3426 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 2,2-DIALKYLGLYCINE DECARBOXYLASE (PYRUVATE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	0	0
			3252	2051	576	607	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	HIS	GLN	CONFLICT	UNP P16932
A	81	GLU	GLY	CONFLICT	UNP P16932

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

- Molecule 4 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 6 is water.

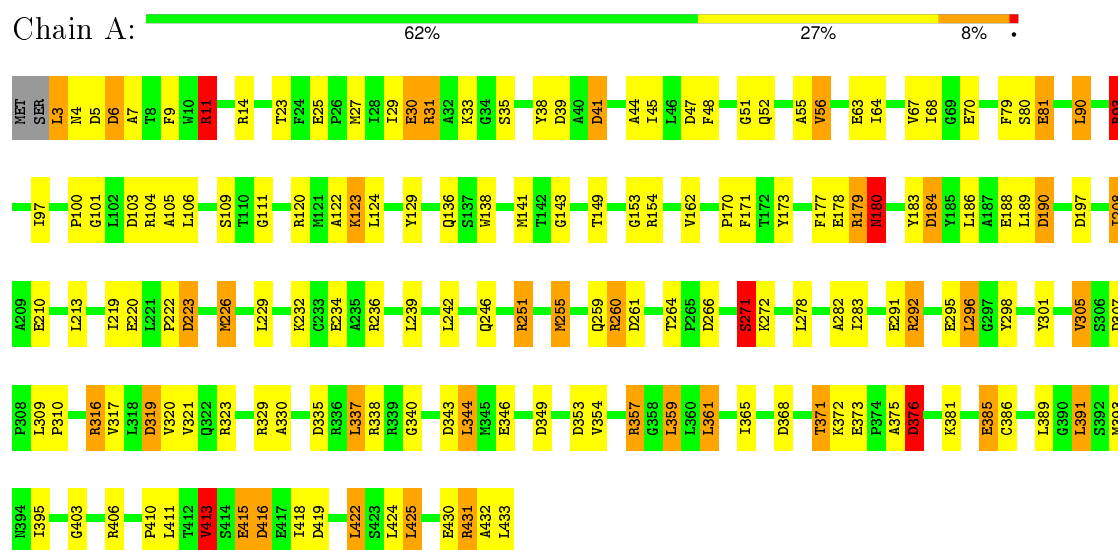
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	145	Total 145	O 145	0	0

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.

Note EDS was not executed.

- Molecule 1: 2,2-DIALKYLGLYCINE DECARBOXYLASE (PYRUVATE)



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	152.70 Å 152.70 Å 86.60 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.176 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3426	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

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: NA, K, MES, PLP

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	0.99	21/3309 (0.6%)	1.41	52/4478 (1.2%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	51	GLY	C-N	9.59	1.56	1.34
1	A	52	GLN	C-N	7.69	1.51	1.34
1	A	220	GLU	CD-OE1	7.37	1.33	1.25
1	A	178	GLU	CD-OE2	6.35	1.32	1.25
1	A	70	GLU	CD-OE2	6.26	1.32	1.25
1	A	234	GLU	CD-OE1	6.16	1.32	1.25
1	A	385	GLU	CD-OE2	5.91	1.32	1.25
1	A	430	GLU	CD-OE2	5.80	1.32	1.25
1	A	346	GLU	CD-OE2	5.78	1.32	1.25
1	A	101	GLY	CA-C	5.72	1.60	1.51
1	A	373	GLU	CD-OE1	5.68	1.31	1.25
1	A	188	GLU	CD-OE2	5.67	1.31	1.25
1	A	295	GLU	CD-OE1	5.58	1.31	1.25
1	A	81	GLU	CD-OE2	5.54	1.31	1.25
1	A	415	GLU	CD-OE2	5.51	1.31	1.25
1	A	210	GLU	CD-OE1	5.44	1.31	1.25
1	A	271	SER	CB-OG	-5.30	1.35	1.42
1	A	63	GLU	CD-OE2	5.22	1.31	1.25
1	A	291	GLU	CD-OE2	5.20	1.31	1.25
1	A	25	GLU	CD-OE1	5.10	1.31	1.25
1	A	30	GLU	CD-OE2	5.08	1.31	1.25

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	6	ASP	CB-CG-OD1	8.70	126.13	118.30
1	A	307	ASP	CB-CG-OD2	-8.66	110.51	118.30
1	A	190	ASP	CB-CG-OD2	-8.44	110.70	118.30
1	A	6	ASP	CB-CG-OD2	-8.21	110.91	118.30
1	A	307	ASP	CB-CG-OD1	7.72	125.25	118.30
1	A	100	PRO	C-N-CA	-7.44	106.68	122.30
1	A	5	ASP	CB-CG-OD2	-7.37	111.66	118.30
1	A	349	ASP	CB-CG-OD2	-7.19	111.83	118.30
1	A	271	SER	N-CA-CB	-7.08	99.88	110.50
1	A	14	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	A	190	ASP	CB-CG-OD1	6.85	124.47	118.30
1	A	266	ASP	CB-CG-OD2	-6.67	112.30	118.30
1	A	197	ASP	CB-CG-OD1	6.63	124.27	118.30
1	A	343	ASP	CB-CG-OD2	-6.59	112.37	118.30
1	A	323	ARG	N-CA-CB	6.53	122.36	110.60
1	A	349	ASP	CB-CG-OD1	6.45	124.11	118.30
1	A	266	ASP	CB-CG-OD1	6.41	124.07	118.30
1	A	11	ARG	CD-NE-CZ	6.39	132.55	123.60
1	A	319	ASP	CB-CG-OD1	6.38	124.04	118.30
1	A	416	ASP	CB-CG-OD1	-6.38	112.56	118.30
1	A	343	ASP	CB-CG-OD1	6.33	124.00	118.30
1	A	368	ASP	CB-CG-OD1	-6.33	112.61	118.30
1	A	416	ASP	CB-CG-OD2	6.29	123.96	118.30
1	A	261	ASP	CB-CG-OD1	-6.25	112.67	118.30
1	A	223	ASP	CB-CG-OD1	-6.21	112.72	118.30
1	A	353	ASP	CB-CG-OD2	-6.19	112.73	118.30
1	A	419	ASP	CB-CG-OD2	-6.15	112.77	118.30
1	A	419	ASP	CB-CG-OD1	6.13	123.82	118.30
1	A	197	ASP	CB-CG-OD2	-6.09	112.82	118.30
1	A	413	VAL	CB-CA-C	-6.09	99.82	111.40
1	A	368	ASP	CB-CG-OD2	6.02	123.72	118.30
1	A	5	ASP	CB-CG-OD1	5.96	123.66	118.30
1	A	251	ARG	NE-CZ-NH1	-5.89	117.35	120.30
1	A	305	VAL	CB-CA-C	-5.84	100.29	111.40
1	A	184	ASP	CB-CG-OD2	-5.78	113.10	118.30
1	A	319	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	A	51	GLY	O-C-N	5.49	131.48	122.70
1	A	103	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	323	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	A	93	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	376	ASP	CB-CG-OD1	-5.32	113.51	118.30
1	A	39	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	260	ARG	NE-CZ-NH2	5.28	122.94	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	ASP	CB-CG-OD1	-5.28	113.55	118.30
1	A	353	ASP	CB-CG-OD1	5.27	123.05	118.30
1	A	47	ASP	CB-CG-OD1	5.26	123.04	118.30
1	A	47	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	A	41	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	A	261	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	14	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	A	406	ARG	NE-CZ-NH2	5.09	122.85	120.30
1	A	9	PHE	CB-CG-CD2	-5.00	117.30	120.80

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	3252	0	3277	82	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	15	0	6	0	0
5	A	12	0	13	1	0
6	A	145	0	0	10	0
All	All	3426	0	3296	83	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 13.

All (83) 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:410:PRO:O	1:A:413:VAL:HG22	1.83	0.79
1:A:338:ARG:HD3	6:A:630:HOH:O	1.85	0.75
1:A:31:ARG:HD3	6:A:609:HOH:O	1.85	0.75
1:A:309:LEU:HB3	1:A:310:PRO:HD3	1.70	0.73
1:A:3:LEU:HD12	1:A:4:ASN:H	1.55	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:TYR:CE1	1:A:44:ALA:HB2	2.29	0.67
1:A:93:ARG:NH2	1:A:319:ASP:OD1	2.27	0.67
1:A:11:ARG:HH11	1:A:11:ARG:HG3	1.59	0.66
1:A:340:GLY:HA3	1:A:422:LEU:HD21	1.78	0.66
1:A:337:LEU:HD13	1:A:361:LEU:HD23	1.78	0.66
1:A:179:ARG:HB2	1:A:184:ASP:CB	2.26	0.65
1:A:179:ARG:HB2	1:A:184:ASP:HB2	1.78	0.65
1:A:271:SER:HB2	6:A:521:HOH:O	1.96	0.65
1:A:64:ILE:O	1:A:68:ILE:HG13	1.98	0.64
1:A:330:ALA:HA	1:A:359:LEU:HD22	1.80	0.63
1:A:329:ARG:HD3	6:A:611:HOH:O	2.00	0.61
1:A:38:TYR:HE1	1:A:44:ALA:HB2	1.66	0.60
1:A:170:PRO:HG2	1:A:222:PRO:HD3	1.84	0.60
1:A:208:ILE:H	1:A:208:ILE:HD12	1.66	0.60
1:A:3:LEU:HD13	1:A:41:ASP:OD2	2.04	0.58
1:A:27:MET:HE1	1:A:29:ILE:HD11	1.87	0.57
1:A:27:MET:CE	1:A:29:ILE:HD11	2.35	0.57
1:A:395:ILE:HD11	1:A:403:GLY:HA3	1.86	0.57
1:A:170:PRO:HG2	1:A:222:PRO:CD	2.36	0.56
1:A:141:MET:HA	1:A:141:MET:CE	2.36	0.56
1:A:226:MET:HE2	1:A:229:LEU:HD23	1.89	0.54
1:A:79:PHE:CE2	1:A:81:GLU:HB2	2.43	0.54
1:A:33:LYS:HG2	6:A:614:HOH:O	2.07	0.54
1:A:305:VAL:HG13	1:A:305:VAL:O	2.09	0.53
1:A:292:ARG:CD	1:A:296:LEU:HD22	2.40	0.51
1:A:317:VAL:O	1:A:321:VAL:HG23	2.11	0.51
1:A:381:LYS:HB3	1:A:432:ALA:HB1	1.94	0.50
1:A:386:CYS:HB3	1:A:391:LEU:HB3	1.94	0.49
1:A:213:LEU:N	1:A:213:LEU:HD12	2.27	0.49
1:A:385:GLU:OE1	1:A:431:ARG:HD2	2.12	0.49
1:A:162:VAL:HA	6:A:616:HOH:O	2.13	0.49
1:A:123:LYS:CD	1:A:129:TYR:HA	2.44	0.48
1:A:393:MET:HB3	6:A:604:HOH:O	2.13	0.47
1:A:93:ARG:O	1:A:97:ILE:HG12	2.14	0.47
1:A:316:ARG:HE	1:A:316:ARG:HB2	1.35	0.47
1:A:173:TYR:OH	1:A:372:LYS:HG2	2.15	0.47
1:A:122:ALA:HB2	1:A:239:LEU:HD12	1.96	0.47
1:A:389:LEU:HB3	1:A:424:LEU:HD22	1.97	0.46
1:A:251:ARG:NE	6:A:510:HOH:O	2.30	0.46
1:A:80:SER:HB2	1:A:301:TYR:CZ	2.50	0.46
1:A:11:ARG:HH11	1:A:11:ARG:CG	2.27	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:PHE:O	1:A:183:TYR:HD1	1.99	0.46
5:A:434:MES:H32	5:A:434:MES:H81	1.46	0.46
1:A:246:GLN:O	1:A:251:ARG:HD2	2.17	0.45
1:A:329:ARG:NH1	6:A:611:HOH:O	2.48	0.45
1:A:123:LYS:HD3	1:A:129:TYR:HA	1.98	0.45
1:A:190:ASP:OD1	1:A:232:LYS:HE3	2.15	0.45
1:A:56:VAL:HG22	1:A:411:LEU:HG	1.97	0.45
1:A:260:ARG:NH1	1:A:357:ARG:CG	2.81	0.44
1:A:171:PHE:CE1	1:A:173:TYR:HB3	2.51	0.44
1:A:141:MET:HE2	1:A:141:MET:HA	1.99	0.44
1:A:242:LEU:HD23	1:A:242:LEU:HA	1.76	0.44
1:A:3:LEU:HD13	1:A:3:LEU:HA	1.82	0.43
1:A:106:LEU:O	1:A:282:ALA:HA	2.18	0.43
1:A:138:TRP:HA	1:A:149:THR:HG23	1.99	0.43
1:A:344:LEU:HA	1:A:344:LEU:HD12	1.74	0.43
1:A:179:ARG:HE	1:A:179:ARG:HB3	1.69	0.43
1:A:255:MET:HB2	1:A:259:GLN:HE22	1.84	0.43
1:A:271:SER:HB3	1:A:272:LYS:H	1.29	0.43
1:A:7:ALA:O	1:A:11:ARG:HB2	2.19	0.43
1:A:90:LEU:HD22	1:A:90:LEU:O	2.19	0.42
1:A:105:ALA:HA	1:A:283:ILE:O	2.19	0.42
1:A:305:VAL:CG1	1:A:305:VAL:O	2.67	0.42
1:A:371:THR:O	1:A:372:LYS:HB2	2.19	0.42
1:A:171:PHE:CZ	1:A:173:TYR:HB3	2.55	0.42
1:A:365:ILE:HG22	1:A:375:ALA:HB3	2.01	0.41
1:A:3:LEU:CD1	1:A:4:ASN:H	2.30	0.41
1:A:27:MET:HE3	1:A:45:ILE:HD11	2.01	0.41
1:A:180:ASN:HA	1:A:180:ASN:HD22	1.69	0.41
1:A:111:GLY:HA2	1:A:271:SER:OG	2.20	0.41
1:A:153:GLY:HA2	6:A:623:HOH:O	2.20	0.41
1:A:236:ARG:HG3	1:A:236:ARG:HH11	1.85	0.41
1:A:359:LEU:HD12	1:A:359:LEU:HA	1.84	0.41
1:A:213:LEU:HB2	1:A:219:ILE:HB	2.01	0.41
1:A:124:LEU:HD22	1:A:298:TYR:HA	2.03	0.41
1:A:425:LEU:HD23	1:A:425:LEU:HA	1.82	0.41
1:A:27:MET:HE1	1:A:45:ILE:HD13	2.03	0.40
1:A:292:ARG:O	1:A:292:ARG:HD3	2.21	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	429/433 (99%)	401 (94%)	22 (5%)	6 (1%)	14 28

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	ALA
1	A	143	GLY
1	A	376	ASP
1	A	6	ASP
1	A	271	SER
1	A	180	ASN

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	334/336 (99%)	285 (85%)	49 (15%)	4 6

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	11	ARG
1	A	23	THR
1	A	30	GLU
1	A	31	ARG

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Mol	Chain	Res	Type
1	A	35	SER
1	A	48	PHE
1	A	56	VAL
1	A	67	VAL
1	A	90	LEU
1	A	93	ARG
1	A	104	ARG
1	A	109	SER
1	A	120	ARG
1	A	123	LYS
1	A	136	GLN
1	A	154	ARG
1	A	179	ARG
1	A	180	ASN
1	A	186	LEU
1	A	189	LEU
1	A	208	ILE
1	A	223	ASP
1	A	226	MET
1	A	255	MET
1	A	264	THR
1	A	271	SER
1	A	278	LEU
1	A	292	ARG
1	A	296	LEU
1	A	316	ARG
1	A	320	VAL
1	A	337	LEU
1	A	344	LEU
1	A	354	VAL
1	A	357	ARG
1	A	359	LEU
1	A	361	LEU
1	A	371	THR
1	A	376	ASP
1	A	391	LEU
1	A	413	VAL
1	A	415	GLU
1	A	416	ASP
1	A	418	ILE
1	A	422	LEU
1	A	425	LEU

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Mol	Chain	Res	Type
1	A	431	ARG
1	A	433	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	180	ASN
1	A	259	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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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$
5	MES	A	434	-	11,12,12	0.53	0	14,16,16	1.03	0
4	PLP	A	437	1	15,15,16	1.25	1 (6%)	21,22,23	2.30	9 (42%)

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
5	MES	A	434	-	-	0/6/14/14	0/1/1/1
4	PLP	A	437	1	-	0/6/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	437	PLP	C4A-C4	-3.04	1.45	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	437	PLP	C5-C6-N1	-3.16	118.37	123.86
4	A	437	PLP	C3-C4-C5	-3.01	115.50	118.78
4	A	437	PLP	C3-C2-N1	-2.64	116.97	120.61
4	A	437	PLP	C5A-C5-C6	-2.44	114.67	119.28
4	A	437	PLP	C4A-C4-C5	2.06	123.03	120.88
4	A	437	PLP	C6-N1-C2	2.79	124.97	119.28
4	A	437	PLP	C2A-C2-C3	3.51	125.27	121.04
4	A	437	PLP	O4P-C5A-C5	4.02	115.64	108.99
4	A	437	PLP	C6-C5-C4	5.27	122.62	118.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	434	MES	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.