Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:52 AM GMT

PDB ID	:	2OKJ
Title	:	The X-ray crystal structure of the 67kDa isoform of Glutamic Acid Decarboxylase (GAD67)
Authors	:	Buckle, A.M.; Fenalti, G.; Law, R.H.P.; Whisstock, J.C.
Deposited on	:	2007-01-17
Resolution	:	2.30 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

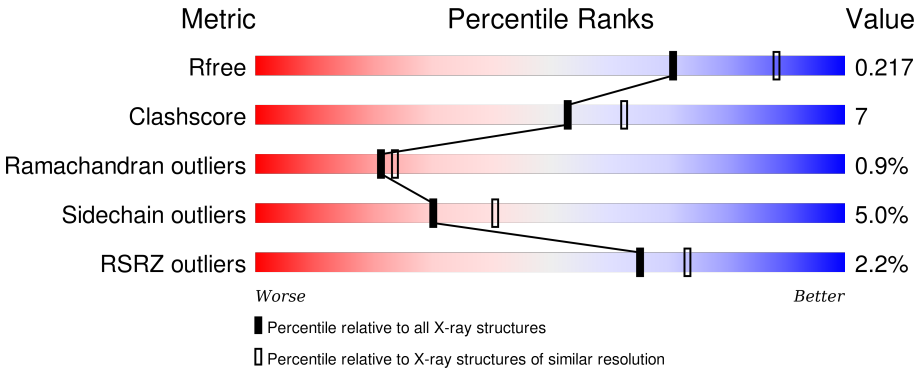
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 <=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	504	3% 87% 10%
1	B	504	2% 88% 10%

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
2	ABU	B	1002[A]	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8314 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.

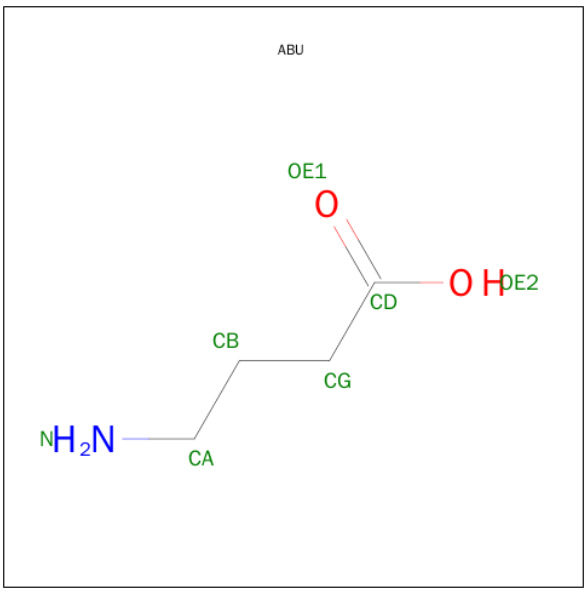
- Molecule 1 is a protein called Glutamate decarboxylase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	P	S	0	0	0
			3947	2535	667	718	1	26			
1	B	504	Total	C	N	O	P	S	0	1	0
			3972	2554	674	717	1	26			

There are 6 discrepancies between the modelled and reference sequences:

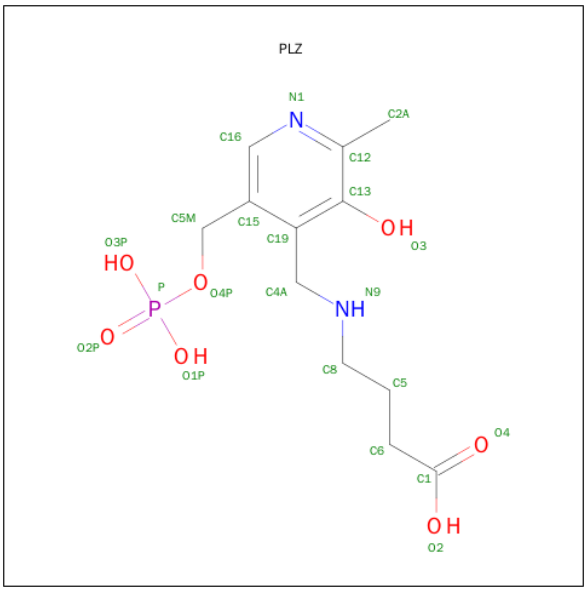
Chain	Residue	Modelled	Actual	Comment	Reference
A	405	LLP	LYS	MODIFIED RESIDUE	UNP Q99259
A	595	HIS	-	EXPRESSION TAG	UNP Q99259
A	596	HIS	-	EXPRESSION TAG	UNP Q99259
B	405	LLP	LYS	MODIFIED RESIDUE	UNP Q99259
B	595	HIS	-	EXPRESSION TAG	UNP Q99259
B	596	HIS	-	EXPRESSION TAG	UNP Q99259

- Molecule 2 is GAMMA-AMINO-BUTANOIC ACID (three-letter code: ABU) (formula: $C_4H_9NO_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 7 4 1 2	0	0
2	B	1	Total C N O 7 4 1 2	0	1

- Molecule 3 is 4-[({3-HYDROXY-2-METHYL-5-[(PHOSPHONOOXY)METHYL]PYRIDIN-4-YL}METHYL)AMINO]BUTANOIC ACID (three-letter code: PLZ) (formula: $C_{12}H_{19}N_2O_7P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C N O P 22 12 2 7 1	0	1

- Molecule 4 is water.

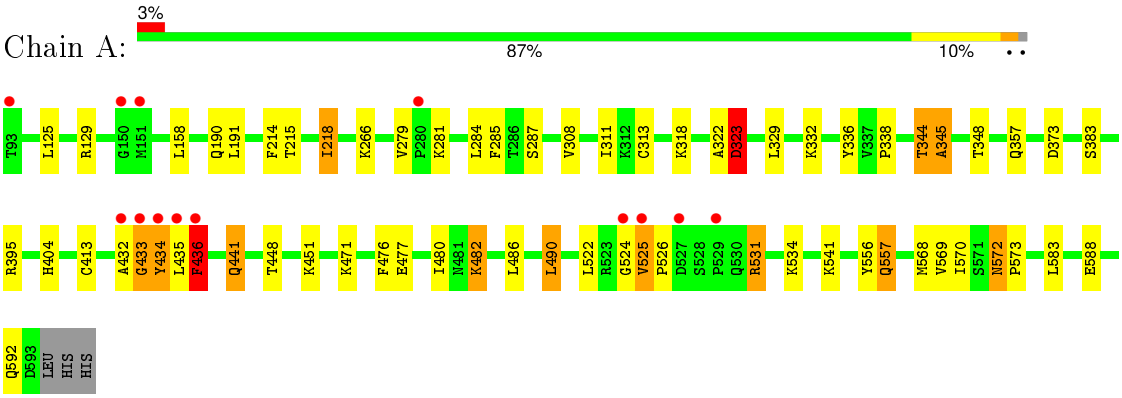
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	177	Total O 177 177	0	0
4	B	182	Total O 182 182	0	0

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.

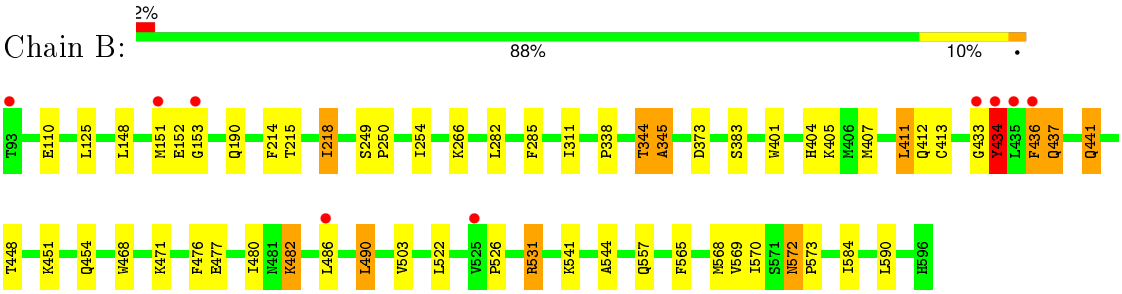
- Molecule 1: Glutamate decarboxylase 1

Chain A:



- Molecule 1: Glutamate decarboxylase 1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	84.05Å 62.74Å 101.35Å 90.00° 106.68° 90.00°	Depositor
Resolution (Å)	97.13 – 2.30 73.14 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.4 (97.13-2.30) 93.4 (73.14-2.30)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.78 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.186 , 0.222 0.183 , 0.217	Depositor DCC
R_{free} test set	2118 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	25.9	Xtriage
Anisotropy	0.273	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 40.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 42271 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8314	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: LLP, ABU, PLZ

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	0.37	0/4015	0.54	1/5428 (0.0%)
1	B	0.36	0/4041	0.51	0/5464
All	All	0.36	0/8056	0.53	1/10892 (0.0%)

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	A	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	323	ASP	N-CA-CB	5.12	119.82	110.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	322	ALA	Peptide

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	3947	0	3866	52	0
1	B	3972	0	3893	51	0
2	A	7	0	5	2	0
2	B	7	0	5	4	0
3	B	22	0	14	5	0
4	A	177	0	0	1	0
4	B	182	0	0	1	0
All	All	8314	0	7783	102	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 7.

All (102) 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:433:GLY:HA3	1:B:434:TYR:CB	1.62	1.26
1:B:433:GLY:CA	1:B:434:TYR:HB3	1.63	1.25
1:A:344:THR:HA	1:A:345:ALA:HB3	1.34	1.09
1:B:344:THR:HA	1:B:345:ALA:HB3	1.34	1.07
1:A:448:THR:HG23	1:A:451:LYS:HD3	1.49	0.92
1:A:433:GLY:HA3	1:A:434:TYR:CD2	2.06	0.90
1:B:344:THR:HA	1:B:345:ALA:CB	2.05	0.85
1:A:344:THR:HA	1:A:345:ALA:CB	2.06	0.85
1:A:490:LEU:HD11	1:A:568:MET:HG2	1.60	0.82
1:A:434:TYR:HA	1:A:435:LEU:HB3	1.63	0.81
1:B:490:LEU:HD11	1:B:568:MET:HG2	1.64	0.80
1:A:482:LYS:HE3	1:A:573:PRO:HA	1.73	0.70
1:B:436:PHE:HA	1:B:437:GLN:HB2	1.77	0.67
1:A:348:THR:HG21	1:B:434:TYR:CE1	2.30	0.67
1:A:348:THR:HG21	1:B:434:TYR:HE1	1.60	0.66
1:A:191:LEU:H	2:A:1001:ABU:HB1	1.62	0.65
1:A:433:GLY:HA3	1:A:434:TYR:HD2	1.59	0.65
1:B:448:THR:HG23	1:B:451:LYS:HD3	1.79	0.65
1:A:404:HIS:CD2	1:A:413:CYS:H	2.16	0.64
1:B:436:PHE:CA	1:B:437:GLN:HB2	2.27	0.63
1:B:344:THR:CA	1:B:345:ALA:HB3	2.22	0.63
1:A:285:PHE:HE1	1:A:338:PRO:HB3	1.63	0.62
1:A:486:LEU:HD12	1:A:570:ILE:HD13	1.83	0.61
1:A:490:LEU:CD1	1:A:568:MET:HG2	2.31	0.60
1:A:215:THR:OG1	1:A:218:ILE:HD12	2.02	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:HIS:CD2	1:B:413:CYS:H	2.20	0.59
1:A:432:ALA:O	1:A:436:PHE:O	2.20	0.59
1:B:250:PRO:HG2	1:B:254:ILE:HD12	1.85	0.58
1:B:522:LEU:O	1:B:531:ARG:NH1	2.36	0.58
1:B:215:THR:OG1	1:B:218:ILE:HD12	2.03	0.58
1:B:486:LEU:HD12	1:B:570:ILE:HD13	1.85	0.57
1:A:279:VAL:O	1:A:279:VAL:HG23	2.04	0.57
1:A:345:ALA:HB3	1:A:373:ASP:O	2.05	0.57
1:B:404:HIS:HD2	1:B:413:CYS:H	1.52	0.57
1:B:345:ALA:HB3	1:B:373:ASP:O	2.05	0.56
1:A:433:GLY:HA3	1:A:434:TYR:CG	2.40	0.56
1:A:434:TYR:HA	1:A:435:LEU:CB	2.31	0.56
1:A:522:LEU:O	1:A:531:ARG:NH1	2.38	0.55
1:A:441:GLN:HG2	1:B:544:ALA:CB	2.37	0.55
1:A:404:HIS:HD2	1:A:413:CYS:H	1.52	0.54
1:A:344:THR:CA	1:A:345:ALA:HB3	2.22	0.54
1:A:344:THR:CA	1:A:345:ALA:CB	2.82	0.53
1:A:433:GLY:HA3	1:A:434:TYR:CB	2.38	0.53
1:B:190:GLN:HE21	2:B:1002[A]:ABU:HB1	1.72	0.53
1:A:285:PHE:HB3	1:A:311:ILE:HD11	1.90	0.52
1:A:435:LEU:O	1:A:435:LEU:HG	2.09	0.52
1:B:572:ASN:HD22	1:B:573:PRO:HD2	1.74	0.52
1:B:436:PHE:HA	1:B:437:GLN:CB	2.40	0.51
1:A:572:ASN:HD22	1:A:573:PRO:HD2	1.76	0.51
1:A:190:GLN:HE21	2:A:1001:ABU:HB2	1.75	0.51
1:B:441:GLN:H	1:B:441:GLN:HE21	1.59	0.50
1:B:476:PHE:O	1:B:480:ILE:HG12	2.10	0.50
1:B:490:LEU:CD1	1:B:568:MET:HG2	2.38	0.50
1:A:476:PHE:O	1:A:480:ILE:HG12	2.12	0.50
1:A:441:GLN:HG2	1:B:544:ALA:HB2	1.94	0.49
1:A:284:LEU:HD13	1:A:308:VAL:HG22	1.95	0.49
1:A:348:THR:CG2	1:B:434:TYR:HE1	2.26	0.49
1:B:285:PHE:HB3	1:B:311:ILE:HD11	1.94	0.49
1:A:345:ALA:CB	1:A:373:ASP:O	2.61	0.48
1:B:383:SER:HA	1:B:477:GLU:HG3	1.95	0.48
1:B:345:ALA:CB	1:B:373:ASP:O	2.60	0.48
1:B:486:LEU:HD11	1:B:570:ILE:HG21	1.95	0.47
1:A:525:VAL:CB	1:A:526:PRO:HA	2.45	0.47
1:A:129:ARG:NH2	1:B:110:GLU:OE2	2.44	0.47
1:B:285:PHE:HE1	1:B:338:PRO:HB3	1.80	0.46
1:A:486:LEU:HD11	1:A:570:ILE:HG21	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:433:GLY:HA3	1:B:434:TYR:HB3	0.70	0.45
1:B:250:PRO:HD3	1:B:454:GLN:NE2	2.31	0.45
1:B:411:LEU:HD23	1:B:412:GLN:HG3	1.98	0.45
1:A:158:LEU:HD22	1:B:468:TRP:HB3	1.99	0.45
1:A:357:GLN:HG3	1:A:395:ARG:HH21	1.82	0.44
1:B:401:TRP:HE1	1:B:407:MET:HE2	1.83	0.44
1:B:249:SER:HB2	1:B:250:PRO:HD2	2.00	0.44
1:A:448:THR:CG2	1:A:451:LYS:HD3	2.36	0.44
1:B:482:LYS:HE2	1:B:573:PRO:HA	2.00	0.44
1:A:557:GLN:CB	1:B:434:TYR:HA	2.49	0.43
1:A:556:TYR:O	1:B:437:GLN:NE2	2.52	0.43
1:B:471:LYS:NZ	4:B:2100:HOH:O	2.51	0.43
1:A:383:SER:HA	1:A:477:GLU:HG3	2.02	0.42
1:B:344:THR:CA	1:B:345:ALA:CB	2.81	0.42
1:A:434:TYR:CG	1:A:434:TYR:O	2.73	0.41
1:B:151:MET:HA	1:B:152:GLU:CB	2.50	0.41
1:A:524:GLY:HA2	1:A:525:VAL:HA	1.70	0.41
1:B:152:GLU:HA	1:B:153:GLY:HA2	1.74	0.41
1:B:285:PHE:CE1	1:B:338:PRO:HB3	2.56	0.41
1:A:281:LYS:HG2	1:A:336:TYR:CE2	2.56	0.41
1:A:287:SER:OG	1:A:313:CYS:SG	2.79	0.41
1:A:525:VAL:CB	1:A:526:PRO:CA	2.99	0.41
1:A:471:LYS:NZ	4:A:1116:HOH:O	2.53	0.40
1:A:313:CYS:HA	1:A:318:LYS:O	2.21	0.40
1:B:490:LEU:HD21	1:B:584:ILE:HD11	2.03	0.40
1:B:503:VAL:HG21	1:B:565:PHE:CG	2.55	0.40
1:B:411:LEU:HA	1:B:412:GLN:HA	1.83	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	498/504 (99%)	470 (94%)	23 (5%)	5 (1%)	19	21
1	B	501/504 (99%)	479 (96%)	18 (4%)	4 (1%)	24	27
All	All	999/1008 (99%)	949 (95%)	41 (4%)	9 (1%)	21	24

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	436	PHE
1	A	433	GLY
1	A	525	VAL
1	B	434	TYR
1	A	323	ASP
1	A	345	ALA
1	B	345	ALA
1	B	437	GLN
1	B	526	PRO

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	409/430 (95%)	387 (95%)	22 (5%)	27	36
1	B	412/430 (96%)	393 (95%)	19 (5%)	33	44
All	All	821/860 (96%)	780 (95%)	41 (5%)	30	41

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	LEU
1	A	214	PHE
1	A	218	ILE
1	A	266	LYS
1	A	323	ASP
1	A	329	LEU
1	A	332	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	344	THR
1	A	434	TYR
1	A	436	PHE
1	A	441	GLN
1	A	482	LYS
1	A	490	LEU
1	A	531	ARG
1	A	534	LYS
1	A	541	LYS
1	A	557	GLN
1	A	569	VAL
1	A	572	ASN
1	A	583	LEU
1	A	588	GLU
1	A	592	GLN
1	B	125	LEU
1	B	148	LEU
1	B	214	PHE
1	B	218	ILE
1	B	266	LYS
1	B	282	LEU
1	B	344	THR
1	B	411	LEU
1	B	434	TYR
1	B	436	PHE
1	B	441	GLN
1	B	482	LYS
1	B	490	LEU
1	B	531	ARG
1	B	541	LYS
1	B	557	GLN
1	B	569	VAL
1	B	572	ASN
1	B	590	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	ASN
1	A	256	ASN
1	A	388	HIS
1	A	404	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	429	GLN
1	A	454	GLN
1	A	458	HIS
1	A	478	ASN
1	A	481	ASN
1	A	564	ASN
1	A	572	ASN
1	A	592	GLN
1	B	142	HIS
1	B	146	GLN
1	B	210	ASN
1	B	256	ASN
1	B	289	GLN
1	B	388	HIS
1	B	404	HIS
1	B	429	GLN
1	B	441	GLN
1	B	454	GLN
1	B	458	HIS
1	B	478	ASN
1	B	481	ASN
1	B	557	GLN
1	B	564	ASN
1	B	572	ASN
1	B	592	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

2 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	A	405	1	23,24,25	1.92	6 (26%)	28,32,34	2.11	7 (25%)
1	LLP	B	405[A]	-	23,24,25	1.92	7 (30%)	28,32,34	2.15	7 (25%)

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	A	405	1	-	0/15/17/19	0/1/1/1
1	LLP	B	405[A]	-	-	0/15/17/19	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	405	LLP	O3-C3	-5.79	1.23	1.37
1	B	405[A]	LLP	O3-C3	-5.69	1.23	1.37
1	A	405	LLP	P-OP2	-3.38	1.42	1.54
1	B	405[A]	LLP	P-OP2	-3.28	1.42	1.54
1	A	405	LLP	P-OP1	-2.58	1.42	1.51
1	B	405[A]	LLP	P-OP1	-2.46	1.43	1.51
1	B	405[A]	LLP	CE-NZ	2.03	1.51	1.46
1	B	405[A]	LLP	C2-N1	2.14	1.38	1.34
1	A	405	LLP	C2-N1	2.15	1.38	1.34
1	A	405	LLP	C4'-NZ	2.31	1.34	1.27
1	B	405[A]	LLP	C4'-NZ	2.45	1.34	1.27
1	A	405	LLP	C4-C4'	2.74	1.51	1.46
1	B	405[A]	LLP	C4-C4'	2.95	1.51	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	405	LLP	OP2-P-OP1	-5.74	92.11	110.58
1	B	405[A]	LLP	OP2-P-OP1	-5.53	92.79	110.58
1	A	405	LLP	C4-C4'-NZ	-2.61	110.53	125.06
1	B	405[A]	LLP	C4-C4'-NZ	-2.57	110.78	125.06
1	B	405[A]	LLP	O-C-CA	-2.42	119.19	125.49
1	A	405	LLP	O-C-CA	-2.39	119.25	125.49
1	A	405	LLP	C5-C6-N1	-2.15	120.14	123.86
1	B	405[A]	LLP	C5-C6-N1	-2.10	120.21	123.86
1	B	405[A]	LLP	OP3-P-OP1	2.53	118.74	110.58
1	A	405	LLP	OP3-P-OP1	2.73	119.38	110.58

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	405	LLP	OP3-P-OP2	3.18	119.48	107.38
1	B	405[A]	LLP	OP3-P-OP2	3.58	121.03	107.38
1	A	405	LLP	OP4-C5'-C5	6.72	120.11	108.99
1	B	405[A]	LLP	OP4-C5'-C5	7.01	120.58	108.99

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
1	B	405[A]	LLP	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

3 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	ABU	A	1001	-	3,6,6	0.21	0	3,6,6	0.44	0
2	ABU	B	1002[A]	3	3,6,6	0.24	0	3,6,6	0.42	0
3	PLZ	B	2001[B]	1,2	19,22,22	3.47	4 (21%)	25,30,30	1.21	1 (4%)

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	ABU	A	1001	-	-	0/2/4/4	0/0/0/0
2	ABU	B	1002[A]	3	-	0/2/4/4	0/0/0/0
3	PLZ	B	2001[B]	1,2	-	0/12/14/14	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2001[B]	PLZ	C4A-C19	-2.85	1.48	1.51
3	B	2001[B]	PLZ	C13-C19	6.02	1.49	1.40
3	B	2001[B]	PLZ	C15-C19	6.36	1.49	1.40
3	B	2001[B]	PLZ	C13-C12	11.80	1.48	1.40

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	2001[B]	PLZ	C16-N1-C12	2.37	124.12	119.28

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	ABU	2	0
2	B	1002[A]	ABU	4	0
3	B	2001[B]	PLZ	5	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, 95^{th} 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	500/504 (99%)	-0.07	13 (2%) 59 68	12, 23, 46, 58	0
1	B	503/504 (99%)	-0.08	9 (1%) 71 78	12, 23, 45, 60	0
All	All	1003/1008 (99%)	-0.08	22 (2%) 65 73	12, 23, 46, 60	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	435	LEU	6.5
1	B	93	THR	6.3
1	A	436	PHE	5.8
1	A	93	THR	5.1
1	A	434	TYR	4.8
1	B	434	TYR	4.5
1	A	433	GLY	4.3
1	A	524	GLY	4.2
1	B	151	MET	3.4
1	B	433	GLY	2.9
1	A	432	ALA	2.9
1	B	525	VAL	2.8
1	A	151	MET	2.8
1	B	486	LEU	2.4
1	A	280	PRO	2.4
1	B	436	PHE	2.3
1	A	150	GLY	2.3
1	A	527	ASP	2.3
1	B	153	GLY	2.2
1	A	525	VAL	2.2
1	A	529	PRO	2.2
1	B	435	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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
1	LLP	A	405	24/25	0.97	0.12	-	12,19,21,24	0
1	LLP	B	405[A]	24/25	0.97	0.12	-	12,19,21,25	15

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, 95^{th} 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
2	ABU	A	1001	7/7	0.84	0.24	0.99	38,39,41,42	0
2	ABU	B	1002[A]	7/7	0.83	0.32	0.73	35,35,37,41	3
3	PLZ	B	2001[B]	22/22	0.96	0.15	-0.42	15,20,40,41	20

6.5 Other polymers (i)

There are no such residues in this entry.