



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:33 AM GMT

PDB ID : 2DGK
Title : Crystal structure of an N-terminal deletion mutant of Escherichia coli GadB
in an autoinhibited state (aldamine)
Authors : Gruetter, M.G.; Capitani, G.; Gut, H.
Deposited on : 2006-03-14
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

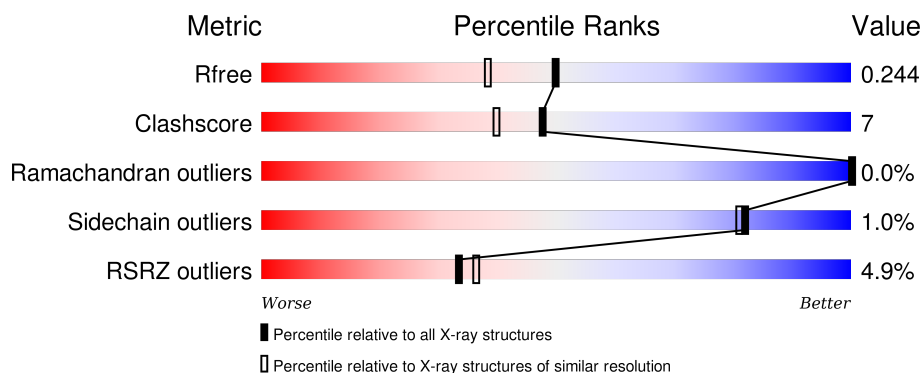
1 Overall quality at a glance

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	452	<div> <div>5%</div> <div>82%</div> <div>14%</div> <div>.</div> </div>
1	B	452	<div> <div>3%</div> <div>78%</div> <div>18%</div> <div>.</div> </div>
1	C	452	<div> <div>4%</div> <div>80%</div> <div>17%</div> <div>.</div> </div>
1	D	452	<div> <div>5%</div> <div>82%</div> <div>14%</div> <div>.</div> </div>
1	E	452	<div> <div>6%</div> <div>78%</div> <div>18%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	452	

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	EDO	B	2191[A]	-	-	-	X
4	EDO	B	2191[B]	-	-	-	X
4	EDO	C	2192[A]	-	-	-	X
4	EDO	C	2192[B]	-	-	-	X
4	EDO	D	2196	-	-	-	X
4	EDO	E	2197	-	-	-	X
4	EDO	F	2198	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 22818 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 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	438	Total	C	N	O	S	0	2	0
			3497	2236	596	640	25			
1	B	438	Total	C	N	O	S	0	3	0
			3497	2234	596	642	25			
1	C	437	Total	C	N	O	S	0	1	0
			3485	2227	595	638	25			
1	D	438	Total	C	N	O	S	0	2	0
			3497	2236	596	640	25			
1	E	438	Total	C	N	O	S	0	3	0
			3499	2236	596	642	25			
1	F	438	Total	C	N	O	S	0	2	0
			3494	2231	597	641	25			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	MET	DELETION	UNP P69910
A	?	-	ASP	DELETION	UNP P69910
A	?	-	LYS	DELETION	UNP P69910
A	?	-	LYS	DELETION	UNP P69910
A	?	-	GLN	DELETION	UNP P69910
A	?	-	VAL	DELETION	UNP P69910
A	?	-	THR	DELETION	UNP P69910
A	?	-	ASP	DELETION	UNP P69910
A	?	-	LEU	DELETION	UNP P69910
A	?	-	ARG	DELETION	UNP P69910
A	?	-	SER	DELETION	UNP P69910
A	?	-	GLU	DELETION	UNP P69910
A	?	-	LEU	DELETION	UNP P69910
A	?	-	LEU	DELETION	UNP P69910
B	?	-	MET	DELETION	UNP P69910
B	?	-	ASP	DELETION	UNP P69910
B	?	-	LYS	DELETION	UNP P69910

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LYS	DELETION	UNP P69910
B	?	-	GLN	DELETION	UNP P69910
B	?	-	VAL	DELETION	UNP P69910
B	?	-	THR	DELETION	UNP P69910
B	?	-	ASP	DELETION	UNP P69910
B	?	-	LEU	DELETION	UNP P69910
B	?	-	ARG	DELETION	UNP P69910
B	?	-	SER	DELETION	UNP P69910
B	?	-	GLU	DELETION	UNP P69910
B	?	-	LEU	DELETION	UNP P69910
B	?	-	LEU	DELETION	UNP P69910
C	?	-	MET	DELETION	UNP P69910
C	?	-	ASP	DELETION	UNP P69910
C	?	-	LYS	DELETION	UNP P69910
C	?	-	LYS	DELETION	UNP P69910
C	?	-	GLN	DELETION	UNP P69910
C	?	-	VAL	DELETION	UNP P69910
C	?	-	THR	DELETION	UNP P69910
C	?	-	ASP	DELETION	UNP P69910
C	?	-	LEU	DELETION	UNP P69910
C	?	-	ARG	DELETION	UNP P69910
C	?	-	SER	DELETION	UNP P69910
C	?	-	GLU	DELETION	UNP P69910
C	?	-	LEU	DELETION	UNP P69910
C	?	-	LEU	DELETION	UNP P69910
D	?	-	MET	DELETION	UNP P69910
D	?	-	ASP	DELETION	UNP P69910
D	?	-	LYS	DELETION	UNP P69910
D	?	-	LYS	DELETION	UNP P69910
D	?	-	GLN	DELETION	UNP P69910
D	?	-	VAL	DELETION	UNP P69910
D	?	-	THR	DELETION	UNP P69910
D	?	-	ASP	DELETION	UNP P69910
D	?	-	LEU	DELETION	UNP P69910
D	?	-	ARG	DELETION	UNP P69910
D	?	-	SER	DELETION	UNP P69910
D	?	-	GLU	DELETION	UNP P69910
D	?	-	LEU	DELETION	UNP P69910
D	?	-	LEU	DELETION	UNP P69910
E	?	-	MET	DELETION	UNP P69910
E	?	-	ASP	DELETION	UNP P69910
E	?	-	LYS	DELETION	UNP P69910

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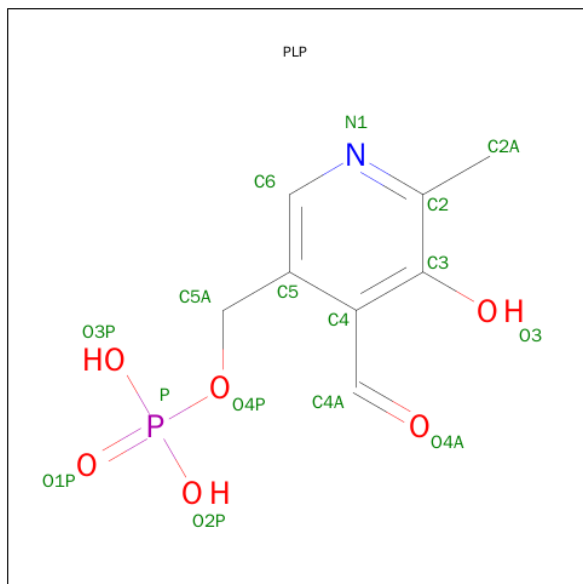
Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	LYS	DELETION	UNP P69910
E	?	-	GLN	DELETION	UNP P69910
E	?	-	VAL	DELETION	UNP P69910
E	?	-	THR	DELETION	UNP P69910
E	?	-	ASP	DELETION	UNP P69910
E	?	-	LEU	DELETION	UNP P69910
E	?	-	ARG	DELETION	UNP P69910
E	?	-	SER	DELETION	UNP P69910
E	?	-	GLU	DELETION	UNP P69910
E	?	-	LEU	DELETION	UNP P69910
E	?	-	LEU	DELETION	UNP P69910
F	?	-	MET	DELETION	UNP P69910
F	?	-	ASP	DELETION	UNP P69910
F	?	-	LYS	DELETION	UNP P69910
F	?	-	LYS	DELETION	UNP P69910
F	?	-	GLN	DELETION	UNP P69910
F	?	-	VAL	DELETION	UNP P69910
F	?	-	THR	DELETION	UNP P69910
F	?	-	ASP	DELETION	UNP P69910
F	?	-	LEU	DELETION	UNP P69910
F	?	-	ARG	DELETION	UNP P69910
F	?	-	SER	DELETION	UNP P69910
F	?	-	GLU	DELETION	UNP P69910
F	?	-	LEU	DELETION	UNP P69910
F	?	-	LEU	DELETION	UNP P69910

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



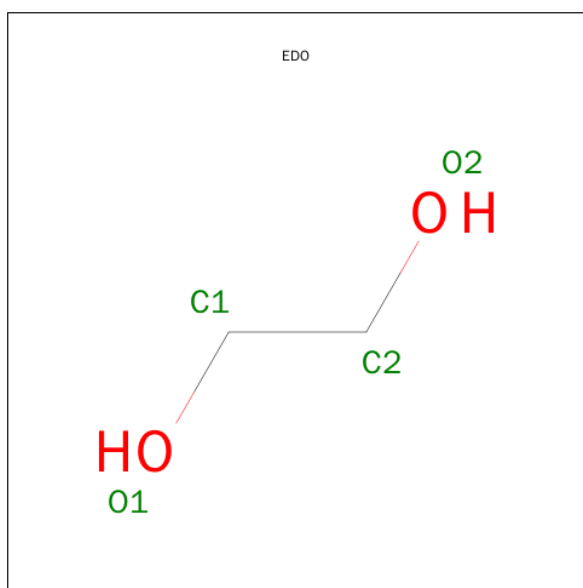
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	E	1	Total O S 5 4 1	0	0
2	F	1	Total O S 5 4 1	0	0

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P 15 8 1 5 1	0	0
3	B	1	Total C N O P 15 8 1 5 1	0	0
3	C	1	Total C N O P 15 8 1 5 1	0	0
3	D	1	Total C N O P 15 8 1 5 1	0	0
3	E	1	Total C N O P 15 8 1 5 1	0	0
3	F	1	Total C N O P 15 8 1 5 1	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 5 2 3	0	1
4	C	1	Total C O 5 2 3	0	1
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	267	Total O 267 267	0	0
5	B	301	Total O 301 301	0	0
5	C	327	Total O 327 327	0	0

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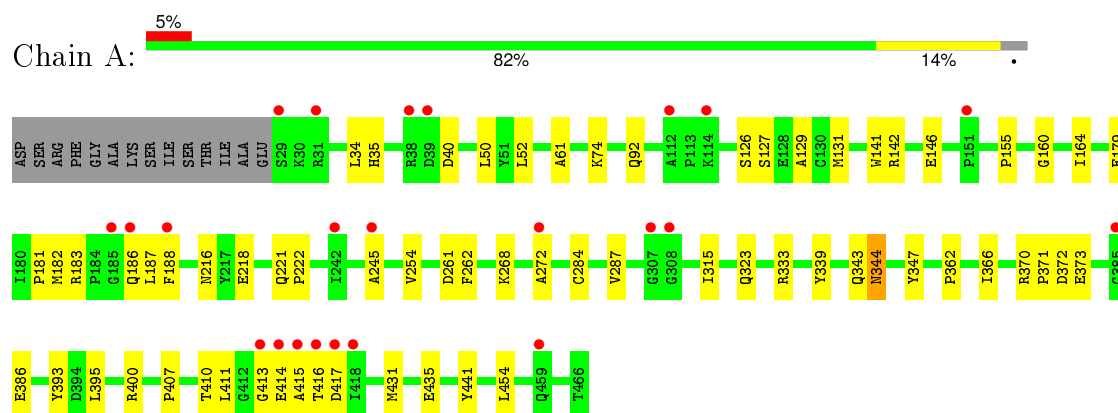
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	311	Total 311	O 311	0	0
5	E	231	Total 231	O 231	0	0
5	F	258	Total 258	O 258	0	0

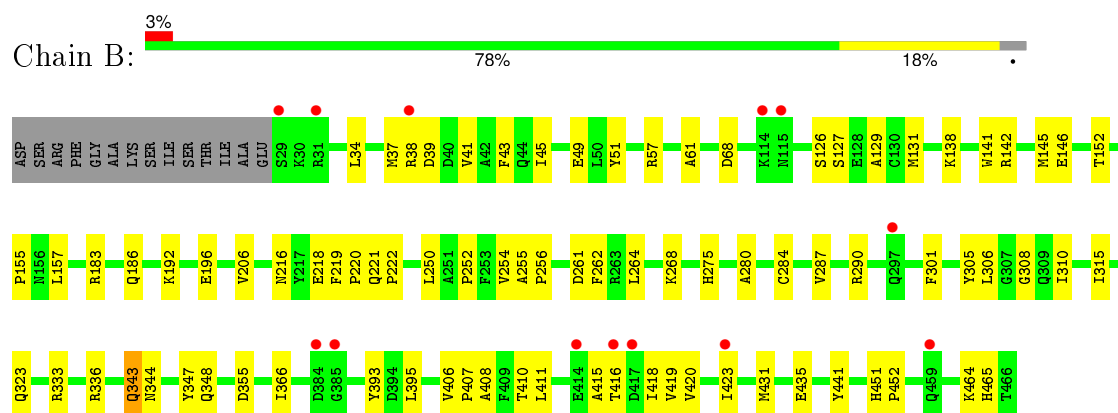
3 Residue-property plots

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.

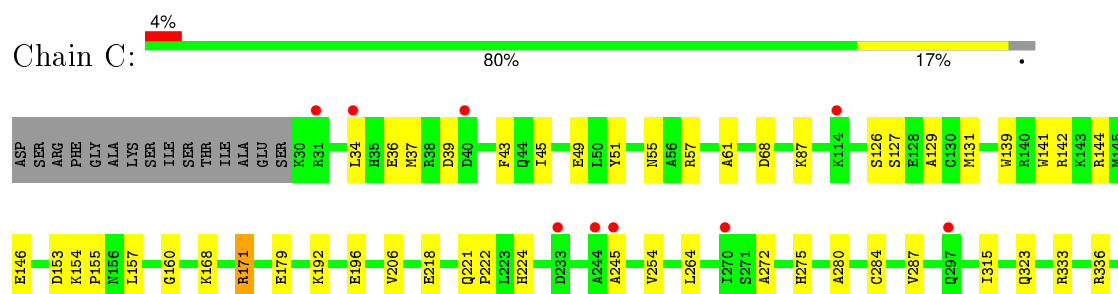
• Molecule 1: Glutamate decarboxylase beta

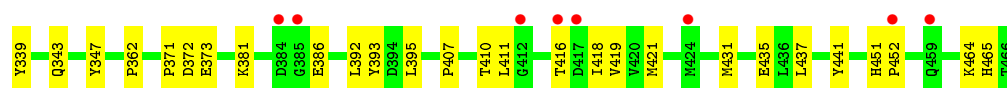


• Molecule 1: Glutamate decarboxylase beta

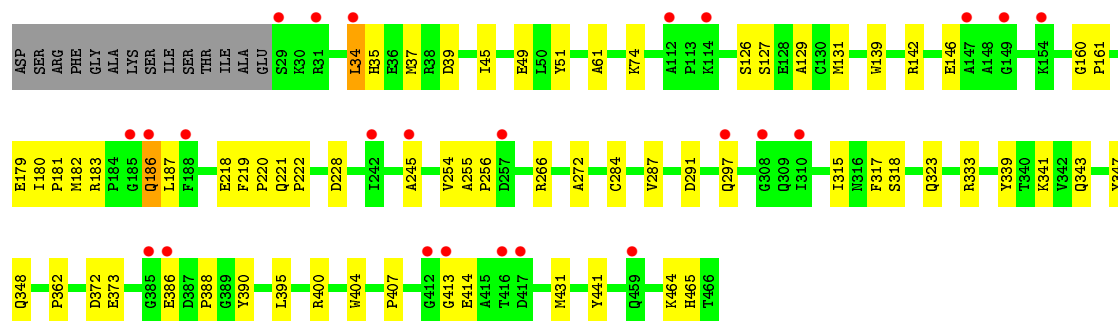
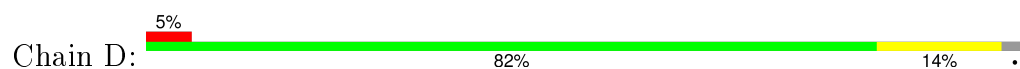


• Molecule 1: Glutamate decarboxylase beta

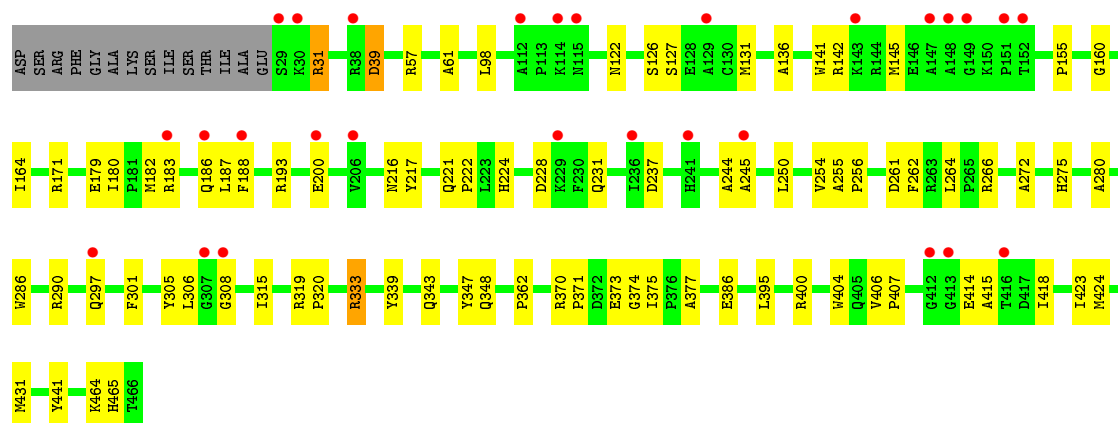
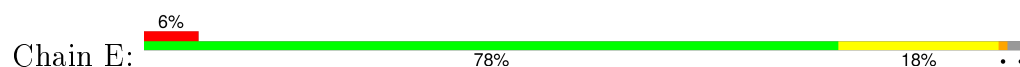




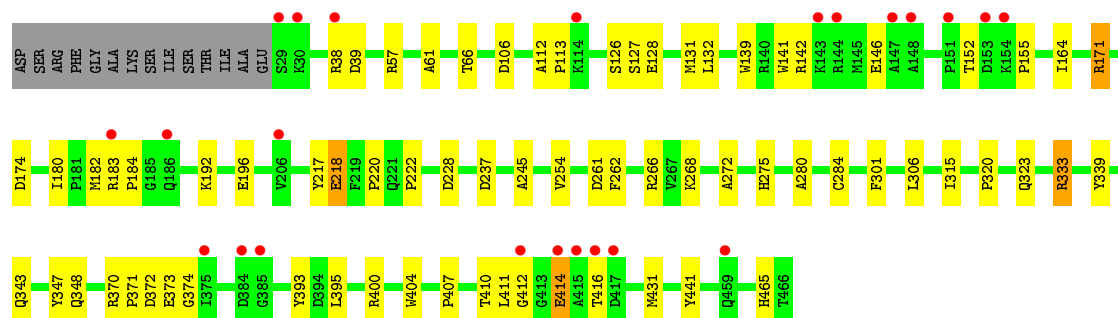
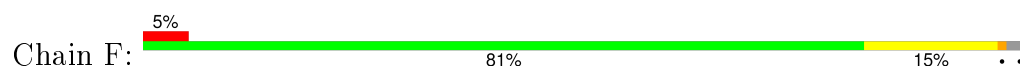
• Molecule 1: Glutamate decarboxylase beta



• Molecule 1: Glutamate decarboxylase beta



• Molecule 1: Glutamate decarboxylase beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.77Å 158.56Å 201.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 29.33 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-1.90) 99.0 (29.33-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.93 (at 1.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.246 0.217 , 0.244	Depositor DCC
R_{free} test set	2308 reflections (1.00%)	DCC
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.3	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	2 of 230716 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	22818	wwPDB-VP
Average B, all atoms (Å ²)	28.0	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 42.27 % 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 2.0904e-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

5.1 Standard geometry

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

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	0.33	0/3601	0.58	1/4881 (0.0%)
1	B	0.33	0/3605	0.59	1/4887 (0.0%)
1	C	0.34	0/3583	0.58	1/4857 (0.0%)
1	D	0.33	0/3601	0.58	1/4881 (0.0%)
1	E	0.31	0/3610	0.57	0/4893
1	F	0.31	0/3598	0.57	1/4878 (0.0%)
All	All	0.33	0/21598	0.58	5/29277 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	218	GLU	N-CA-C	-5.36	96.54	111.00
1	D	218	GLU	N-CA-C	-5.35	96.57	111.00
1	F	218	GLU	N-CA-C	-5.34	96.57	111.00
1	A	218	GLU	N-CA-C	-5.22	96.92	111.00
1	C	218	GLU	N-CA-C	-5.14	97.12	111.00

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	3497	0	3394	47	0
1	B	3497	0	3396	60	0
1	C	3485	0	3385	59	0
1	D	3497	0	3395	52	0
1	E	3499	0	3388	73	0
1	F	3494	0	3389	56	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	1	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	15	0	6	1	0
3	B	15	0	6	1	0
3	C	15	0	6	1	0
3	D	15	0	6	1	0
3	E	15	0	6	1	0
3	F	15	0	6	1	0
4	A	4	0	6	0	0
4	B	9	0	12	0	0
4	C	9	0	12	0	0
4	D	4	0	6	0	0
4	E	4	0	6	0	0
4	F	4	0	6	0	0
5	A	267	0	0	2	0
5	B	301	0	0	3	0
5	C	327	0	0	1	0
5	D	311	0	0	3	0
5	E	231	0	0	5	0
5	F	258	0	0	4	0
All	All	22818	0	20431	306	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 (306) 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:250:LEU:HD23	1:B:343:GLN:HG3	1.56	0.86
1:A:362:PRO:HB2	1:A:386:GLU:HG2	1.60	0.83
1:C:221:GLN:HB3	1:C:222:PRO:HD3	1.61	0.83
1:D:362:PRO:HB2	1:D:386:GLU:HG2	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:183:ARG:HG2	1:E:414:GLU:HB2	1.61	0.82
1:E:221:GLN:HB3	1:E:222:PRO:HD3	1.59	0.82
1:A:333:ARG:HD3	1:B:39:ASP:HA	1.63	0.81
1:C:333:ARG:HD3	1:D:39:ASP:HA	1.63	0.81
1:A:221:GLN:HB3	1:A:222:PRO:HD3	1.64	0.79
1:D:221:GLN:HB3	1:D:222:PRO:HD3	1.63	0.79
1:E:362:PRO:HB2	1:E:386:GLU:HG2	1.63	0.79
1:E:127:SER:O	1:E:131[B]:MET:HG3	1.83	0.78
1:C:127:SER:O	1:C:131[B]:MET:HG2	1.83	0.77
1:E:127:SER:O	1:E:131[A]:MET:HG2	1.84	0.77
1:C:39:ASP:HA	1:D:333:ARG:HD3	1.67	0.76
1:C:362:PRO:HB2	1:C:386:GLU:HG2	1.69	0.74
1:B:127:SER:O	1:B:131[B]:MET:HG2	1.86	0.74
1:F:348:GLN:HG2	1:F:431:MET:HE1	1.69	0.74
1:F:127:SER:O	1:F:131[B]:MET:HG2	1.89	0.72
1:E:228:ASP:HA	1:E:266:ARG:HH21	1.55	0.71
1:A:127:SER:O	1:A:131[B]:MET:HG2	1.92	0.70
1:F:182:MET:HE3	1:F:412:GLY:H	1.57	0.68
1:D:182:MET:H	1:D:413:GLY:HA3	1.61	0.66
1:C:410:THR:HG22	1:C:419:VAL:HG22	1.78	0.66
1:A:181:PRO:HB2	1:A:414:GLU:HG3	1.76	0.66
1:D:142:ARG:O	1:D:146:GLU:HG3	1.95	0.65
1:B:410:THR:HG22	1:B:419:VAL:HG22	1.78	0.65
1:A:362:PRO:CB	1:A:386:GLU:HG2	2.26	0.65
1:C:142:ARG:O	1:C:146:GLU:HG3	1.97	0.65
1:F:171:ARG:HB2	1:F:171:ARG:HH11	1.63	0.64
1:A:431:MET:O	1:A:435:GLU:HG2	1.98	0.64
1:E:348:GLN:HE21	1:E:431:MET:CE	2.11	0.63
1:A:61:ALA:HB2	1:A:407:PRO:HD3	1.80	0.63
1:C:315:ILE:HG23	1:D:131[A]:MET:SD	2.39	0.63
1:D:127:SER:O	1:D:131[B]:MET:HG2	1.98	0.62
1:F:182:MET:CE	1:F:412:GLY:H	2.13	0.61
1:E:348:GLN:HE21	1:E:431:MET:HE3	1.64	0.61
1:F:370:ARG:HD3	1:F:373:GLU:OE2	2.00	0.61
1:C:154:LYS:N	1:C:155:PRO:HD3	2.16	0.61
1:C:171:ARG:HB2	1:C:171:ARG:HH11	1.67	0.60
1:D:362:PRO:CB	1:D:386:GLU:HG2	2.32	0.60
1:A:142:ARG:O	1:A:146:GLU:HG3	2.02	0.59
1:C:381:LYS:HE3	1:C:418:ILE:HD12	1.84	0.59
1:E:31:ARG:HB3	1:E:31:ARG:NH1	2.18	0.59
1:F:61:ALA:HB2	1:F:407:PRO:HD3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:ALA:HB2	1:C:407:PRO:HD3	1.85	0.59
1:C:411:LEU:O	1:C:416:THR:HA	2.03	0.59
1:F:339:TYR:O	1:F:343:GLN:HG2	2.03	0.58
1:A:141:TRP:CZ3	1:A:155:PRO:HB3	2.38	0.58
1:B:57:ARG:NH2	1:B:68:ASP:OD2	2.36	0.58
1:A:216:ASN:HD21	1:A:366:ILE:HG22	1.68	0.58
1:B:250:LEU:HD23	1:B:343:GLN:CG	2.31	0.58
1:E:333:ARG:NE	1:F:39:ASP:OD1	2.35	0.58
1:E:31:ARG:HH11	1:E:31:ARG:HB3	1.69	0.57
1:D:339:TYR:O	1:D:343:GLN:HG2	2.02	0.57
1:C:171:ARG:NH1	1:C:171:ARG:HB2	2.19	0.57
1:B:51:TYR:CB	1:C:57:ARG:HG3	2.34	0.57
1:C:144:ARG:HH11	1:C:144:ARG:HB3	1.69	0.57
1:E:164:ILE:HD11	1:F:301:PHE:HB3	1.87	0.57
1:C:339:TYR:O	1:C:343:GLN:HG2	2.05	0.57
1:B:250:LEU:CD2	1:B:343:GLN:HG3	2.31	0.57
1:A:372:ASP:OD1	1:A:373:GLU:HG3	2.05	0.57
1:B:142:ARG:O	1:B:146:GLU:HG3	2.05	0.57
1:C:315:ILE:HG21	1:D:315:ILE:HD13	1.87	0.56
1:E:39:ASP:HA	1:F:333:ARG:HG3	1.86	0.56
1:D:181:PRO:HB3	1:D:414:GLU:HG3	1.88	0.56
1:A:34:LEU:HD12	1:A:35:HIS:CE1	2.41	0.56
1:C:395:LEU:HD21	1:C:441:TYR:CZ	2.40	0.56
1:B:395:LEU:HD21	1:B:441:TYR:CZ	2.40	0.56
1:E:126:SER:HB2	3:E:1504:PLP:O4P	2.06	0.56
1:D:254:VAL:HG21	1:D:347:TYR:CE2	2.41	0.56
1:E:231:GLN:NE2	1:E:237:ASP:HB2	2.20	0.55
1:C:57:ARG:NH2	1:C:68:ASP:OD2	2.39	0.55
1:C:141:TRP:CZ3	1:C:155:PRO:HG3	2.40	0.55
1:E:315:ILE:HG21	1:F:315:ILE:HD13	1.88	0.55
1:B:305:TYR:HB2	1:B:308:GLY:O	2.06	0.55
1:B:57:ARG:HG3	1:C:51:TYR:CB	2.37	0.55
1:C:393:TYR:OH	1:C:410:THR:HG23	2.07	0.55
1:A:131[A]:MET:SD	1:B:315:ILE:HG23	2.47	0.55
1:C:431:MET:O	1:C:435:GLU:HG3	2.07	0.55
1:D:388:PRO:HB2	1:D:390:TYR:CE2	2.42	0.54
1:A:333:ARG:HE	1:B:39:ASP:HB3	1.73	0.54
1:B:157[B]:LEU:HD11	1:B:206:VAL:HG23	1.89	0.54
1:D:372:ASP:OD1	1:D:373:GLU:HG3	2.07	0.54
1:D:61:ALA:HB2	1:D:407:PRO:HD3	1.90	0.54
1:E:415:ALA:HB1	1:E:418:ILE:HD12	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:TYR:O	1:A:343:GLN:HG2	2.07	0.54
1:A:415:ALA:C	1:A:417:ASP:H	2.11	0.53
1:B:415:ALA:HB1	1:B:418:ILE:HD12	1.89	0.53
1:E:39:ASP:HB2	5:E:4987:HOH:O	2.08	0.53
1:E:145:MET:CE	1:E:200:GLU:HG2	2.39	0.53
1:F:126:SER:HB2	3:F:1505:PLP:O4P	2.08	0.53
1:E:224:HIS:CD2	1:E:264:LEU:HB3	2.44	0.53
1:E:370:ARG:HD3	1:E:373:GLU:OE2	2.08	0.53
1:A:315:ILE:HG23	1:B:131[A]:MET:SD	2.48	0.52
1:E:31:ARG:NH1	1:F:106:ASP:OD2	2.42	0.52
1:B:264:LEU:O	1:B:290:ARG:NH2	2.42	0.52
1:F:393:TYR:OH	1:F:410:THR:HG23	2.09	0.52
1:E:377:ALA:HB2	1:E:424:MET:HE1	1.91	0.52
1:A:254:VAL:HG21	1:A:347:TYR:CE2	2.45	0.52
1:A:74:LYS:HE2	1:B:43:PHE:CZ	2.44	0.52
1:C:157:LEU:HD11	1:C:206:VAL:HG23	1.91	0.52
1:E:301:PHE:HB3	1:F:164:ILE:HD11	1.91	0.51
1:B:221:GLN:HB3	1:B:222:PRO:HD3	1.92	0.51
1:E:255:ALA:N	1:E:256:PRO:HD3	2.24	0.51
1:A:245:ALA:HA	1:A:272:ALA:HA	1.92	0.51
1:B:275:HIS:HA	1:B:280:ALA:O	2.11	0.51
1:E:142:ARG:HH22	1:F:174:ASP:CG	2.13	0.51
1:C:333:ARG:HE	1:D:39:ASP:HB3	1.76	0.51
1:C:131[A]:MET:SD	1:D:315:ILE:HG23	2.51	0.51
1:D:228:ASP:HA	1:D:266:ARG:HH21	1.75	0.51
1:E:160:GLY:O	1:E:179:GLU:HG3	2.11	0.51
1:E:362:PRO:CB	1:E:386:GLU:HG2	2.36	0.51
1:A:411:LEU:O	1:A:416:THR:HA	2.11	0.51
1:B:61:ALA:HB2	1:B:407:PRO:HD3	1.93	0.50
1:C:144:ARG:NH1	1:C:144:ARG:HB3	2.25	0.50
1:C:336:ARG:HD2	1:D:34:LEU:HD22	1.92	0.50
1:E:306:LEU:HD21	1:F:400:ARG:NE	2.25	0.50
1:E:315:ILE:HD13	1:F:315:ILE:HG21	1.93	0.50
1:C:192:LYS:O	1:C:196:GLU:HG3	2.12	0.50
1:A:182:MET:H	1:A:413:GLY:HA3	1.77	0.50
1:C:37:MET:O	1:D:333:ARG:HD2	2.12	0.49
1:E:395:LEU:HD21	1:E:441:TYR:CZ	2.46	0.49
1:F:141:TRP:CZ3	1:F:155:PRO:HB3	2.47	0.49
1:B:343:GLN:OE1	1:B:343:GLN:HA	2.12	0.49
1:E:375:ILE:HD11	1:E:424:MET:HE1	1.93	0.49
1:E:61:ALA:HB2	1:E:407:PRO:HD3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:171:ARG:HH11	1:E:171:ARG:HG2	1.77	0.49
1:E:371:PRO:HD2	5:E:4999:HOH:O	2.13	0.49
1:C:153:ASP:C	1:C:155:PRO:HD3	2.33	0.49
1:C:372:ASP:OD1	1:C:373:GLU:HG3	2.12	0.49
1:B:344:ASN:O	1:B:348:GLN:HG3	2.12	0.49
1:E:136:ALA:HB3	5:E:4801:HOH:O	2.13	0.49
1:E:188[A]:PHE:CZ	1:E:216:ASN:HB2	2.48	0.49
1:B:38:ARG:HB2	1:B:41:VAL:HG23	1.95	0.49
1:F:411:LEU:O	1:F:416:THR:HA	2.12	0.49
1:F:228:ASP:HA	1:F:266:ARG:HH21	1.78	0.48
1:C:392:LEU:HD13	1:C:421:MET:HB2	1.95	0.48
1:F:254:VAL:HG21	1:F:347:TYR:CE2	2.47	0.48
1:B:192:LYS:O	1:B:196:GLU:HG3	2.14	0.48
1:D:51:TYR:CB	1:E:57:ARG:HG3	2.44	0.48
1:A:284:CYS:HB2	1:A:323:GLN:HB3	1.96	0.48
1:D:464:LYS:HE3	5:D:4990:HOH:O	2.13	0.48
1:C:254:VAL:HG21	1:C:347:TYR:CE2	2.48	0.48
1:C:34:LEU:C	1:C:34:LEU:HD23	2.34	0.48
1:A:187:LEU:O	1:A:188[A]:PHE:CD1	2.67	0.48
1:E:131[B]:MET:SD	1:F:315:ILE:HG23	2.54	0.48
1:C:34:LEU:O	1:C:34:LEU:HD23	2.13	0.48
1:D:255:ALA:N	1:D:256:PRO:HD3	2.29	0.48
1:B:431:MET:O	1:B:435:GLU:HG3	2.14	0.47
1:E:264:LEU:O	1:E:290:ARG:NH2	2.48	0.47
1:A:395:LEU:HD21	1:A:441:TYR:CZ	2.49	0.47
1:A:216:ASN:ND2	1:A:366:ILE:HG22	2.29	0.47
1:B:411:LEU:O	1:B:416:THR:HA	2.14	0.47
1:D:182:MET:HG2	1:D:187:LEU:CD2	2.45	0.47
1:A:181:PRO:CB	1:A:414:GLU:HG3	2.42	0.47
1:E:339:TYR:O	1:E:343:GLN:HG2	2.15	0.47
1:A:40:ASP:HB2	5:A:4952:HOH:O	2.15	0.47
1:B:465:HIS:O	1:B:465:HIS:CG	2.68	0.47
1:B:255:ALA:N	1:B:256:PRO:HD3	2.30	0.46
1:D:400:ARG:HA	1:D:404:TRP:O	2.16	0.46
1:A:371:PRO:HD2	5:A:4916:HOH:O	2.15	0.46
1:D:183:ARG:NH2	1:D:186:GLN:OE1	2.45	0.46
1:D:297:GLN:HG2	5:D:4967:HOH:O	2.14	0.46
1:F:465:HIS:CG	1:F:465:HIS:O	2.68	0.46
1:E:186:GLN:HE21	1:E:193:ARG:CZ	2.29	0.46
1:B:138:LYS:HE2	1:B:142:ARG:HH21	1.80	0.46
1:A:126:SER:HB2	3:A:1500:PLP:O4P	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:245:ALA:HA	1:D:272:ALA:HA	1.97	0.46
1:C:275:HIS:HA	1:C:280:ALA:O	2.16	0.46
1:D:160:GLY:O	1:D:179:GLU:HG3	2.16	0.46
1:C:129:ALA:HB1	1:C:287:VAL:HB	1.97	0.46
1:E:306:LEU:HD21	1:F:400:ARG:CZ	2.46	0.46
1:C:284:CYS:HB2	1:C:323:GLN:HB3	1.98	0.46
1:E:400:ARG:HA	1:E:404:TRP:O	2.15	0.46
1:E:141:TRP:CZ3	1:E:155:PRO:HB3	2.50	0.46
1:E:180:ILE:N	1:E:180:ILE:HD12	2.30	0.46
1:E:343:GLN:OE1	1:E:343:GLN:HA	2.16	0.45
1:E:254:VAL:HG21	1:E:347:TYR:CE2	2.51	0.45
1:F:142:ARG:O	1:F:146:GLU:HG3	2.16	0.45
1:E:122:ASN:HB2	1:E:286:TRP:CZ3	2.52	0.45
1:D:129:ALA:HB1	1:D:287:VAL:HB	1.97	0.45
1:F:395:LEU:HD21	1:F:441:TYR:CZ	2.51	0.45
1:A:344:ASN:HA	1:A:344:ASN:HD22	1.60	0.45
1:F:372:ASP:OD1	1:F:373:GLU:HG3	2.17	0.45
1:C:43:PHE:CZ	1:D:74:LYS:HE2	2.51	0.45
1:B:216:ASN:HD21	1:B:366:ILE:HG22	1.82	0.45
1:E:182:MET:HE2	1:E:187:LEU:O	2.16	0.45
1:C:333:ARG:HD3	1:D:39:ASP:CA	2.42	0.45
1:C:55:ASN:O	1:C:57:ARG:NH1	2.50	0.45
1:B:284:CYS:HB2	1:B:323:GLN:HB3	1.98	0.45
1:F:171:ARG:CB	1:F:171:ARG:HH11	2.28	0.45
1:E:145:MET:HE1	1:E:200:GLU:HG2	1.99	0.45
1:D:395:LEU:HD21	1:D:441:TYR:CZ	2.51	0.45
1:E:39:ASP:OD1	1:F:333:ARG:NH2	2.50	0.45
1:F:192:LYS:HG2	1:F:196:GLU:OE2	2.17	0.45
1:C:371:PRO:HD2	5:C:4829:HOH:O	2.17	0.44
1:E:217:TYR:CD2	1:E:374:GLY:HA2	2.51	0.44
1:F:371:PRO:HD2	5:F:4931:HOH:O	2.16	0.44
1:C:245:ALA:HA	1:C:272:ALA:HA	2.00	0.44
1:E:377:ALA:HB2	1:E:424:MET:CE	2.48	0.44
1:F:57:ARG:CZ	1:F:66:THR:O	2.66	0.44
1:A:92:GLN:HG2	1:B:49:GLU:O	2.18	0.44
1:C:224:HIS:CD2	1:C:264:LEU:HB3	2.52	0.44
1:D:465:HIS:O	1:D:465:HIS:CG	2.71	0.44
1:A:164:ILE:HD11	1:B:301:PHE:HB3	1.99	0.44
1:B:183:ARG:NH2	1:B:186:GLN:OE1	2.47	0.44
1:F:180:ILE:HD12	1:F:180:ILE:N	2.32	0.44
1:D:284:CYS:HB2	1:D:323:GLN:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:ARG:HG3	1:C:51:TYR:CG	2.53	0.43
1:F:217:TYR:CD2	1:F:374:GLY:HA2	2.53	0.43
1:B:393:TYR:OH	1:B:410:THR:HG23	2.18	0.43
1:B:41:VAL:O	1:B:45:ILE:HG13	2.18	0.43
1:E:261:ASP:HB2	1:E:262:PHE:H	1.62	0.43
1:F:261:ASP:HB2	1:F:262:PHE:H	1.64	0.43
1:F:142:ARG:HG3	1:F:152:THR:HG21	1.99	0.43
1:C:464:LYS:O	1:C:465:HIS:HB3	2.17	0.43
1:A:400:ARG:NE	1:B:306:LEU:HD21	2.33	0.43
1:C:333:ARG:HD2	1:D:37:MET:O	2.18	0.43
1:A:34:LEU:O	1:A:34:LEU:HD13	2.18	0.43
1:B:141:TRP:CZ3	1:B:155:PRO:HB3	2.53	0.43
1:B:254:VAL:HG21	1:B:347:TYR:CE2	2.53	0.43
1:B:186:GLN:HG3	1:B:186:GLN:O	2.18	0.43
1:B:408:ALA:HA	1:B:420:VAL:O	2.18	0.43
1:E:400:ARG:NH1	1:F:306:LEU:HD11	2.33	0.43
1:C:465:HIS:CG	1:C:465:HIS:O	2.72	0.43
1:D:291:ASP:HB2	2:D:4773:SO4:O3	2.19	0.43
1:C:160:GLY:O	1:C:179:GLU:HG3	2.19	0.43
1:F:183:ARG:NH2	1:F:414:GLU:OE2	2.46	0.43
1:C:45:ILE:O	1:C:49:GLU:HG3	2.19	0.43
1:B:126:SER:HB2	3:B:1501:PLP:O4P	2.18	0.43
1:B:310:ILE:HD11	5:B:5052:HOH:O	2.19	0.43
1:E:465:HIS:O	1:E:465:HIS:CG	2.72	0.43
1:D:182:MET:HG2	1:D:187:LEU:HD22	2.01	0.42
1:C:36:GLU:OE2	1:D:341:LYS:HD3	2.19	0.42
1:C:126:SER:HB2	3:C:1502:PLP:O4P	2.18	0.42
1:F:128:GLU:O	1:F:132:LEU:HG	2.19	0.42
1:D:34:LEU:HD12	1:D:35:HIS:CE1	2.53	0.42
1:D:348:GLN:HG2	1:D:431:MET:HE1	2.00	0.42
1:B:145:MET:HB2	1:B:152:THR:HG22	2.02	0.42
1:A:393:TYR:OH	1:A:410:THR:HG23	2.19	0.42
1:F:171:ARG:CG	1:F:171:ARG:HH11	2.33	0.42
1:E:250:LEU:HD12	1:E:375:ILE:HG22	2.00	0.42
1:A:333:ARG:HD2	1:B:37:MET:O	2.19	0.42
1:F:400:ARG:HA	1:F:404:TRP:O	2.20	0.42
1:A:370:ARG:HA	1:A:371:PRO:HD3	1.92	0.42
1:A:160:GLY:O	1:A:179:GLU:HG3	2.19	0.42
1:D:126:SER:HB2	3:D:1503:PLP:O4P	2.20	0.42
1:E:244:ALA:O	1:E:245:ALA:C	2.58	0.42
1:C:315:ILE:HG21	1:D:315:ILE:CD1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:ALA:C	1:A:417:ASP:N	2.74	0.42
1:A:183:ARG:NH2	1:A:186:GLN:OE1	2.53	0.42
1:C:451:HIS:N	1:C:452:PRO:HD3	2.35	0.42
1:D:45:ILE:O	1:D:49:GLU:HG3	2.20	0.42
1:E:250:LEU:CD1	1:E:375:ILE:HG22	2.50	0.42
1:E:464:LYS:O	1:E:465:HIS:HB3	2.20	0.42
1:F:183:ARG:HG3	1:F:184:PRO:HD2	2.02	0.41
1:D:348:GLN:HG2	1:D:431:MET:CE	2.50	0.41
1:A:261:ASP:HB2	1:A:262:PHE:H	1.64	0.41
1:E:406:VAL:HG13	1:E:423:ILE:HG12	2.02	0.41
1:F:284:CYS:HB2	1:F:323:GLN:HB3	2.02	0.41
1:A:129:ALA:HB1	1:A:287:VAL:HB	2.01	0.41
1:E:98:LEU:HG	5:E:4984:HOH:O	2.20	0.41
1:E:319:ARG:HB2	1:E:320:PRO:HD2	2.01	0.41
1:F:38:ARG:HG3	5:F:4914:HOH:O	2.19	0.41
1:F:275:HIS:HA	1:F:280:ALA:O	2.21	0.41
1:F:320:PRO:HB2	5:F:4978:HOH:O	2.19	0.41
1:D:362:PRO:HG3	1:D:388:PRO:HG3	2.02	0.41
1:B:406:VAL:HG13	1:B:423:ILE:HG12	2.02	0.41
1:E:348:GLN:HG2	1:E:431:MET:HE3	2.02	0.41
1:C:168:LYS:HE2	5:D:4986:HOH:O	2.21	0.41
1:E:221:GLN:HB3	1:E:222:PRO:CD	2.41	0.41
1:D:180:ILE:HA	1:D:181:PRO:HD3	1.92	0.41
1:F:245:ALA:HA	1:F:272:ALA:HA	2.03	0.41
1:B:415:ALA:HA	5:B:4941:HOH:O	2.21	0.41
1:D:362:PRO:CG	1:D:388:PRO:HG3	2.51	0.41
1:E:375:ILE:CD1	1:E:424:MET:HE1	2.51	0.41
1:F:182:MET:HE3	1:F:412:GLY:N	2.32	0.41
1:F:237:ASP:OD1	1:F:266:ARG:NH1	2.52	0.41
1:D:317:PHE:HB3	1:D:318:SER:H	1.77	0.41
1:B:261:ASP:HB2	1:B:262:PHE:H	1.66	0.41
1:B:219:PHE:HA	1:B:220:PRO:HD3	1.90	0.41
1:E:221:GLN:CB	1:E:222:PRO:HD3	2.42	0.41
1:C:362:PRO:CB	1:C:386:GLU:HG2	2.45	0.41
1:F:411:LEU:O	1:F:416:THR:HG22	2.21	0.41
1:B:464:LYS:O	1:B:465:HIS:HB3	2.19	0.41
1:B:216:ASN:ND2	1:B:366:ILE:HG22	2.35	0.41
1:E:333:ARG:HG3	1:F:39:ASP:HA	2.02	0.40
1:E:245:ALA:HA	1:E:272:ALA:HA	2.02	0.40
1:D:219:PHE:HA	1:D:220:PRO:HD3	1.88	0.40
1:A:50:LEU:HD23	1:A:50:LEU:HA	1.96	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:218:GLU:O	1:F:220:PRO:HD3	2.22	0.40
1:B:129:ALA:HB1	1:B:287:VAL:HB	2.03	0.40
1:E:297:GLN:HG2	5:E:4898:HOH:O	2.21	0.40
1:B:451:HIS:N	1:B:452:PRO:HD3	2.36	0.40
1:F:112:ALA:HA	1:F:113:PRO:HD3	1.85	0.40
1:D:161:PRO:HG3	1:D:182:MET:HE3	2.04	0.40
1:A:34:LEU:HD22	1:B:336:ARG:HD2	2.03	0.40
1:E:305:TYR:HB2	1:E:308:GLY:O	2.20	0.40
1:B:252:PRO:HA	5:B:4991:HOH:O	2.22	0.40
1:A:454:LEU:HD21	5:F:4819:HOH:O	2.22	0.40
1:E:275:HIS:HA	1:E:280:ALA:O	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	438/452 (97%)	424 (97%)	14 (3%)	0	100	100
1	B	439/452 (97%)	428 (98%)	11 (2%)	0	100	100
1	C	436/452 (96%)	423 (97%)	13 (3%)	0	100	100
1	D	438/452 (97%)	422 (96%)	15 (3%)	1 (0%)	52	42
1	E	439/452 (97%)	420 (96%)	19 (4%)	0	100	100
1	F	438/452 (97%)	427 (98%)	11 (2%)	0	100	100
All	All	2628/2712 (97%)	2544 (97%)	83 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	186	GLN

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	367/376 (98%)	364 (99%)	3 (1%)	86	86
1	B	368/376 (98%)	363 (99%)	5 (1%)	74	71
1	C	365/376 (97%)	361 (99%)	4 (1%)	80	79
1	D	367/376 (98%)	365 (100%)	2 (0%)	92	92
1	E	368/376 (98%)	365 (99%)	3 (1%)	86	86
1	F	367/376 (98%)	362 (99%)	5 (1%)	74	71
All	All	2202/2256 (98%)	2180 (99%)	22 (1%)	82	81

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	268	LYS
1	A	344	ASN
1	B	34	LEU
1	B	268	LYS
1	B	333	ARG
1	B	343	GLN
1	B	355	ASP
1	C	87	LYS
1	C	139	TRP
1	C	171	ARG
1	C	437	LEU
1	D	34	LEU
1	D	139	TRP
1	E	31	ARG
1	E	39	ASP
1	E	333	ARG
1	F	139	TRP
1	F	171	ARG
1	F	268	LYS
1	F	333	ARG
1	F	414	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	109	HIS
1	A	216	ASN
1	A	309	GLN
1	A	344	ASN
1	A	451	HIS
1	A	459	GLN
1	B	109	HIS
1	B	216	ASN
1	C	109	HIS
1	C	309	GLN
1	C	344	ASN
1	D	309	GLN
1	D	348	GLN
1	E	109	HIS
1	E	186	GLN
1	E	309	GLN
1	E	344	ASN
1	E	348	GLN
1	E	455	GLN
1	F	201	ASN
1	F	309	GLN
1	F	344	ASN
1	F	348	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

22 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$
3	PLP	A	1500	1	15,15,16	1.17	2 (13%)	21,22,23	1.23	5 (23%)
4	EDO	A	2193	-	3,3,3	0.44	0	2,2,2	0.44	0
2	SO4	A	4770	-	4,4,4	1.50	0	6,6,6	1.62	1 (16%)
3	PLP	B	1501	1	15,15,16	0.90	0	21,22,23	1.24	4 (19%)
4	EDO	B	2191[A]	-	3,3,3	0.49	0	2,2,2	0.40	0
4	EDO	B	2191[B]	-	3,3,3	0.49	0	2,2,2	0.39	0
4	EDO	B	2194	-	3,3,3	0.51	0	2,2,2	0.40	0
2	SO4	B	4771	-	4,4,4	1.55	0	6,6,6	1.61	1 (16%)
3	PLP	C	1502	1	15,15,16	0.99	1 (6%)	21,22,23	1.25	5 (23%)
4	EDO	C	2192[A]	-	3,3,3	0.47	0	2,2,2	0.46	0
4	EDO	C	2192[B]	-	3,3,3	0.47	0	2,2,2	0.45	0
4	EDO	C	2195	-	3,3,3	0.46	0	2,2,2	0.45	0
2	SO4	C	4772	-	4,4,4	1.51	0	6,6,6	1.62	1 (16%)
3	PLP	D	1503	1	15,15,16	1.06	1 (6%)	21,22,23	1.23	5 (23%)
4	EDO	D	2196	-	3,3,3	0.48	0	2,2,2	0.44	0
2	SO4	D	4773	-	4,4,4	1.58	0	6,6,6	1.62	1 (16%)
3	PLP	E	1504	1	15,15,16	1.19	1 (6%)	21,22,23	1.25	5 (23%)
4	EDO	E	2197	-	3,3,3	0.43	0	2,2,2	0.47	0
2	SO4	E	4774	-	4,4,4	1.56	0	6,6,6	1.62	1 (16%)
3	PLP	F	1505	1	15,15,16	1.33	2 (13%)	21,22,23	1.20	4 (19%)
4	EDO	F	2198	-	3,3,3	0.43	0	2,2,2	0.48	0
2	SO4	F	4775	-	4,4,4	1.60	0	6,6,6	1.62	1 (16%)

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
3	PLP	A	1500	1	-	0/6/6/8	0/1/1/1
4	EDO	A	2193	-	-	0/1/1/1	0/0/0/0
2	SO4	A	4770	-	-	0/0/0/0	0/0/0/0
3	PLP	B	1501	1	-	0/6/6/8	0/1/1/1
4	EDO	B	2191[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	2191[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	B	2194	-	-	0/1/1/1	0/0/0/0
2	SO4	B	4771	-	-	0/0/0/0	0/0/0/0
3	PLP	C	1502	1	-	0/6/6/8	0/1/1/1
4	EDO	C	2192[A]	-	-	0/1/1/1	0/0/0/0
4	EDO	C	2192[B]	-	-	0/1/1/1	0/0/0/0
4	EDO	C	2195	-	-	0/1/1/1	0/0/0/0
2	SO4	C	4772	-	-	0/0/0/0	0/0/0/0
3	PLP	D	1503	1	-	0/6/6/8	0/1/1/1
4	EDO	D	2196	-	-	0/1/1/1	0/0/0/0
2	SO4	D	4773	-	-	0/0/0/0	0/0/0/0
3	PLP	E	1504	1	-	0/6/6/8	0/1/1/1
4	EDO	E	2197	-	-	0/1/1/1	0/0/0/0
2	SO4	E	4774	-	-	0/0/0/0	0/0/0/0
3	PLP	F	1505	1	-	0/6/6/8	0/1/1/1
4	EDO	F	2198	-	-	0/1/1/1	0/0/0/0
2	SO4	F	4775	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1500	PLP	C5-C4	2.12	1.43	1.40
3	C	1502	PLP	C5-C4	2.15	1.43	1.40
3	D	1503	PLP	C3-C2	2.31	1.42	1.40
3	F	1505	PLP	C5-C4	2.50	1.43	1.40
3	A	1500	PLP	C3-C2	2.81	1.42	1.40
3	E	1504	PLP	C3-C2	3.02	1.42	1.40
3	F	1505	PLP	C3-C2	3.28	1.43	1.40

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1501	PLP	C3-C2-N1	-2.23	117.53	120.61
3	C	1502	PLP	C3-C2-N1	-2.21	117.56	120.61
3	A	1500	PLP	C3-C2-N1	-2.18	117.60	120.61
3	D	1503	PLP	C3-C2-N1	-2.16	117.63	120.61
3	E	1504	PLP	C3-C2-N1	-2.14	117.66	120.61
3	F	1505	PLP	C3-C2-N1	-2.02	117.82	120.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1502	PLP	C2A-C2-C3	2.01	123.46	121.04
3	F	1505	PLP	O3-C3-C2	2.03	121.18	117.66
3	D	1503	PLP	O3-C3-C2	2.07	121.26	117.66
3	D	1503	PLP	C2A-C2-C3	2.13	123.61	121.04
3	C	1502	PLP	O3-C3-C2	2.14	121.38	117.66
3	A	1500	PLP	O3-C3-C2	2.14	121.39	117.66
3	B	1501	PLP	O3-C3-C2	2.15	121.40	117.66
3	E	1504	PLP	O3-C3-C2	2.17	121.43	117.66
3	E	1504	PLP	C2A-C2-C3	2.18	123.66	121.04
3	A	1500	PLP	C2A-C2-C3	2.20	123.69	121.04
3	F	1505	PLP	C6-N1-C2	2.24	123.84	119.28
3	A	1500	PLP	C6-N1-C2	2.24	123.84	119.28
3	E	1504	PLP	C6-N1-C2	2.26	123.89	119.28
3	D	1503	PLP	C6-N1-C2	2.27	123.90	119.28
3	C	1502	PLP	C6-N1-C2	2.32	124.01	119.28
3	B	1501	PLP	C6-N1-C2	2.34	124.06	119.28
3	B	1501	PLP	O3P-P-O1P	2.35	118.14	110.58
3	F	1505	PLP	O3P-P-O1P	2.40	118.32	110.58
3	A	1500	PLP	O3P-P-O1P	2.43	118.41	110.58
3	D	1503	PLP	O3P-P-O1P	2.45	118.47	110.58
3	C	1502	PLP	O3P-P-O1P	2.53	118.73	110.58
3	E	1504	PLP	O3P-P-O1P	2.54	118.75	110.58
2	B	4771	SO4	O4-S-O3	3.76	124.26	108.98
2	D	4773	SO4	O4-S-O3	3.78	124.34	108.98
2	A	4770	SO4	O4-S-O3	3.78	124.36	108.98
2	C	4772	SO4	O4-S-O3	3.78	124.36	108.98
2	E	4774	SO4	O4-S-O3	3.78	124.37	108.98
2	F	4775	SO4	O4-S-O3	3.79	124.40	108.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1500	PLP	1	0
3	B	1501	PLP	1	0
3	C	1502	PLP	1	0
3	D	1503	PLP	1	0
2	D	4773	SO4	1	0
3	E	1504	PLP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1505	PLP	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

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	438/452 (96%)	0.37	23 (5%) 30 33	17, 26, 37, 41	0
1	B	438/452 (96%)	0.31	13 (2%) 54 57	15, 26, 37, 42	0
1	C	437/452 (96%)	0.33	17 (3%) 43 47	15, 26, 37, 42	0
1	D	438/452 (96%)	0.32	24 (5%) 29 32	16, 25, 36, 41	0
1	E	438/452 (96%)	0.56	28 (6%) 23 25	17, 29, 47, 52	0
1	F	438/452 (96%)	0.49	23 (5%) 30 33	16, 29, 43, 48	0
All	All	2627/2712 (96%)	0.40	128 (4%) 33 36	15, 27, 40, 52	0

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	236	ILE	5.2
1	E	114	LYS	5.2
1	E	148	ALA	4.9
1	B	114	LYS	4.2
1	F	114	LYS	4.2
1	A	416	THR	4.1
1	C	385	GLY	4.1
1	E	151	PRO	4.1
1	D	416	THR	4.0
1	A	114	LYS	3.9
1	E	416	THR	3.8
1	F	385	GLY	3.8
1	F	417	ASP	3.8
1	D	149	GLY	3.7
1	F	459	GLN	3.7
1	F	414	GLU	3.7
1	F	153	ASP	3.6
1	C	114	LYS	3.6
1	E	188[A]	PHE	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	114	LYS	3.6
1	B	31	ARG	3.6
1	F	412	GLY	3.4
1	A	188[A]	PHE	3.4
1	F	416	THR	3.4
1	E	29	SER	3.3
1	E	308	GLY	3.2
1	A	31	ARG	3.1
1	D	31	ARG	3.1
1	A	39	ASP	3.1
1	D	147	ALA	3.1
1	E	115	ASN	3.1
1	F	144	ARG	3.0
1	D	112	ALA	3.0
1	E	183	ARG	3.0
1	F	183	ARG	3.0
1	F	29	SER	3.0
1	D	29	SER	2.9
1	D	413	GLY	2.9
1	E	30	LYS	2.8
1	F	143	LYS	2.8
1	B	414	GLU	2.7
1	C	417	ASP	2.7
1	D	188[A]	PHE	2.7
1	A	112	ALA	2.7
1	A	186	GLN	2.7
1	F	148	ALA	2.7
1	E	149	GLY	2.7
1	B	385	GLY	2.7
1	E	297	GLN	2.7
1	D	417	ASP	2.6
1	E	112	ALA	2.6
1	F	415	ALA	2.6
1	B	29	SER	2.6
1	A	29	SER	2.6
1	D	308	GLY	2.6
1	E	152	THR	2.6
1	A	308	GLY	2.6
1	C	459	GLN	2.6
1	B	423	ILE	2.6
1	E	206	VAL	2.6
1	D	459	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	F	30	LYS	2.5
1	E	412	GLY	2.5
1	E	147	ALA	2.5
1	B	115	ASN	2.5
1	C	416	THR	2.5
1	A	185	GLY	2.5
1	D	245	ALA	2.5
1	C	384	ASP	2.5
1	A	245	ALA	2.5
1	B	416	THR	2.5
1	A	459	GLN	2.5
1	D	297	GLN	2.5
1	A	417	ASP	2.5
1	C	233	ASP	2.5
1	A	272	ALA	2.5
1	C	245	ALA	2.4
1	C	40	ASP	2.4
1	C	31	ARG	2.4
1	C	244	ALA	2.4
1	D	385	GLY	2.4
1	B	384	ASP	2.4
1	A	385	GLY	2.3
1	C	412	GLY	2.3
1	A	38	ARG	2.3
1	D	386	GLU	2.3
1	C	270	ILE	2.3
1	B	417	ASP	2.3
1	A	307	GLY	2.3
1	A	414	GLU	2.3
1	B	297	GLN	2.3
1	B	459	GLN	2.3
1	E	241	HIS	2.3
1	F	154	LYS	2.3
1	F	186	GLN	2.3
1	E	307	GLY	2.2
1	E	129	ALA	2.2
1	F	38	ARG	2.2
1	E	413	GLY	2.2
1	A	151	PRO	2.2
1	B	38	ARG	2.2
1	C	34	LEU	2.2
1	D	154	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	229	LYS	2.2
1	F	147	ALA	2.2
1	A	242	ILE	2.2
1	A	413	GLY	2.2
1	D	412	GLY	2.2
1	F	151	PRO	2.2
1	C	424	MET	2.2
1	A	418	ILE	2.1
1	E	200	GLU	2.1
1	D	186	GLN	2.1
1	E	38	ARG	2.1
1	D	257	ASP	2.1
1	C	297	GLN	2.1
1	D	34	LEU	2.1
1	C	452	PRO	2.1
1	E	186	GLN	2.1
1	E	143	LYS	2.1
1	D	310	ILE	2.1
1	A	415	ALA	2.1
1	F	375	ILE	2.0
1	F	206	VAL	2.0
1	F	384	ASP	2.0
1	E	245	ALA	2.0
1	D	185	GLY	2.0
1	D	242	ILE	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	EDO	B	2191[A]	4/4	0.56	0.27	5.32	47,47,47,47	1
4	EDO	B	2191[B]	4/4	0.56	0.27	5.32	47,47,47,47	1
4	EDO	C	2192[B]	4/4	0.62	0.27	4.80	47,47,47,47	1
4	EDO	C	2192[A]	4/4	0.62	0.27	4.80	47,47,47,47	1
4	EDO	E	2197	4/4	0.83	0.19	3.16	38,38,38,39	0
4	EDO	D	2196	4/4	0.92	0.15	2.65	33,33,33,33	0
4	EDO	F	2198	4/4	0.83	0.16	2.15	32,32,32,33	0
4	EDO	A	2193	4/4	0.87	0.16	1.92	32,33,33,33	0
3	PLP	F	1505	15/16	0.95	0.15	1.12	22,23,24,24	0
3	PLP	D	1503	15/16	0.98	0.14	0.42	17,18,19,19	0
3	PLP	E	1504	15/16	0.96	0.14	0.28	22,22,22,22	0
4	EDO	C	2195	4/4	0.90	0.12	0.26	29,30,30,30	0
2	SO4	B	4771	5/5	0.94	0.20	0.04	50,50,50,50	0
3	PLP	C	1502	15/16	0.97	0.13	0.03	18,20,21,21	0
3	PLP	B	1501	15/16	0.96	0.12	0.00	19,20,21,22	0
3	PLP	A	1500	15/16	0.97	0.13	-0.08	19,20,21,21	0
2	SO4	A	4770	5/5	0.90	0.19	-0.46	37,37,37,37	5
2	SO4	C	4772	5/5	0.93	0.14	-0.47	38,38,38,38	5
2	SO4	D	4773	5/5	0.95	0.16	-0.62	31,31,31,31	5
2	SO4	F	4775	5/5	0.91	0.16	-0.71	52,52,52,52	5
2	SO4	E	4774	5/5	0.96	0.14	-0.82	53,53,53,53	5
4	EDO	B	2194	4/4	0.97	0.07	-1.76	27,27,27,27	0

6.5 Other polymers ⓘ

There are no such residues in this entry.