



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:34 AM UTC

PDB ID : 1PMM / pdb_00001pmm
Title : Crystal structure of Escherichia coli GadB (low pH)
Authors : Capitani, G.; De Biase, D.; Aurizi, C.; Gut, H.; Bossa, F.; Grutter, M.G.
Deposited on : 2003-06-11
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

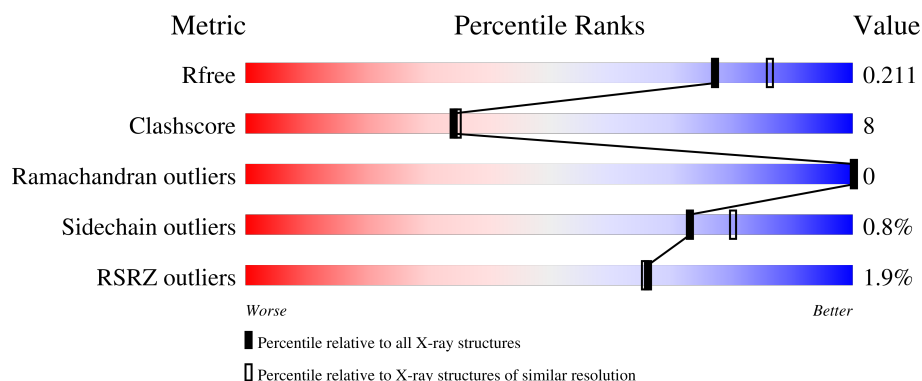
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

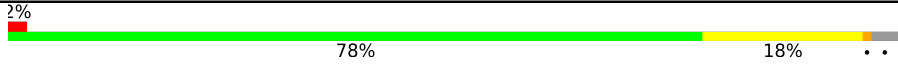
The reported resolution of this entry is 2.00 Å.

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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	466	
1	B	466	
1	C	466	
1	D	466	
1	E	466	

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Mol	Chain	Length	Quality of chain
1	F	466	2% 77% 19%

2 Entry composition [i](#)

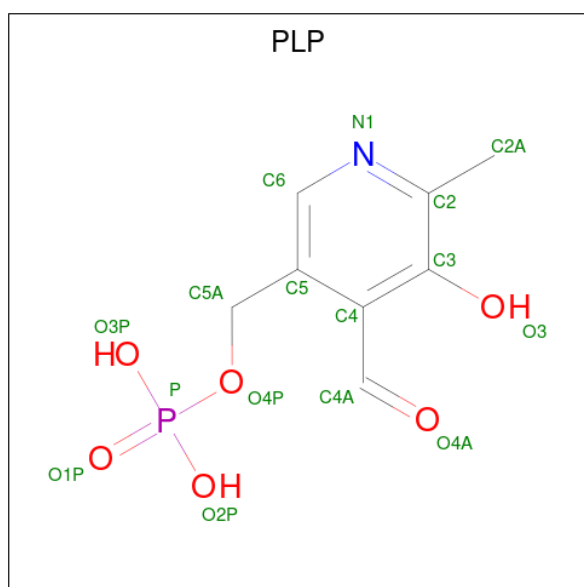
There are 4 unique types of molecules in this entry. The entry contains 23447 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

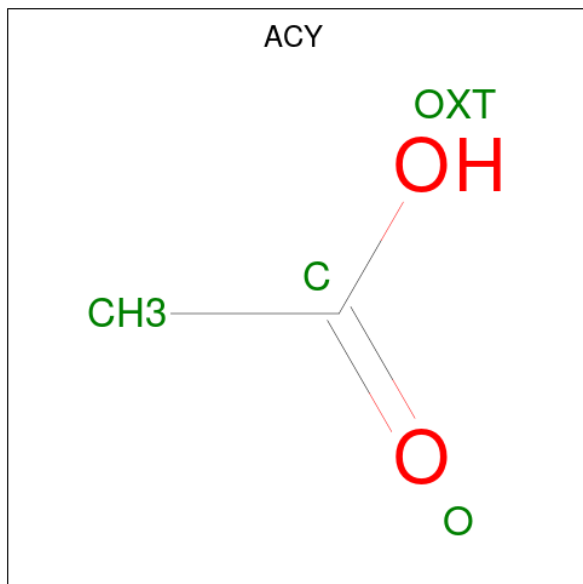
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	3	0
			3584	2287	610	661	26			
1	B	450	Total	C	N	O	S	0	3	0
			3585	2287	611	662	25			
1	C	449	Total	C	N	O	S	0	4	0
			3579	2282	608	664	25			
1	D	450	Total	C	N	O	S	0	3	0
			3584	2287	610	661	26			
1	E	449	Total	C	N	O	S	0	3	0
			3575	2281	608	660	26			
1	F	450	Total	C	N	O	S	0	3	0
			3584	2287	610	661	26			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).



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

- Molecule 3 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

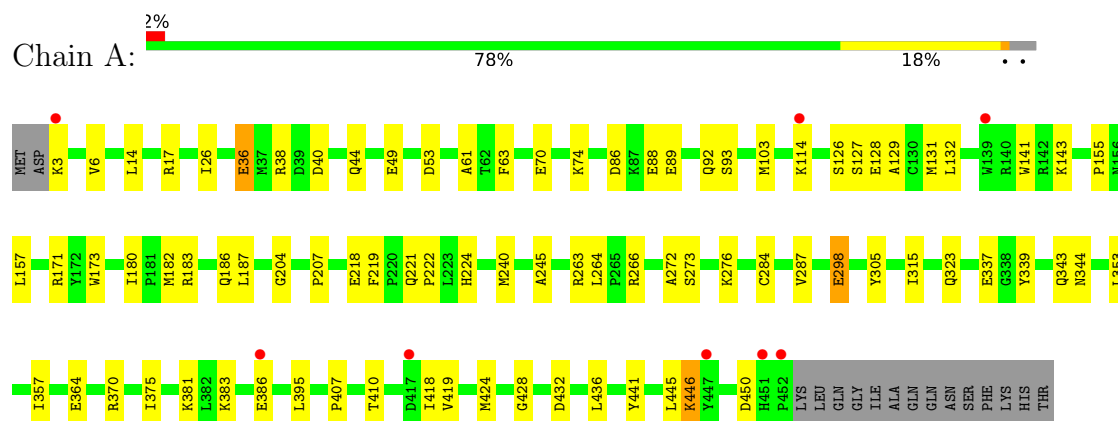
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	309	Total 309	O 309	0	0
4	B	320	Total 320	O 320	0	0
4	C	284	Total 284	O 284	0	0
4	D	333	Total 333	O 333	0	0
4	E	292	Total 292	O 292	0	0
4	F	304	Total 304	O 304	0	0

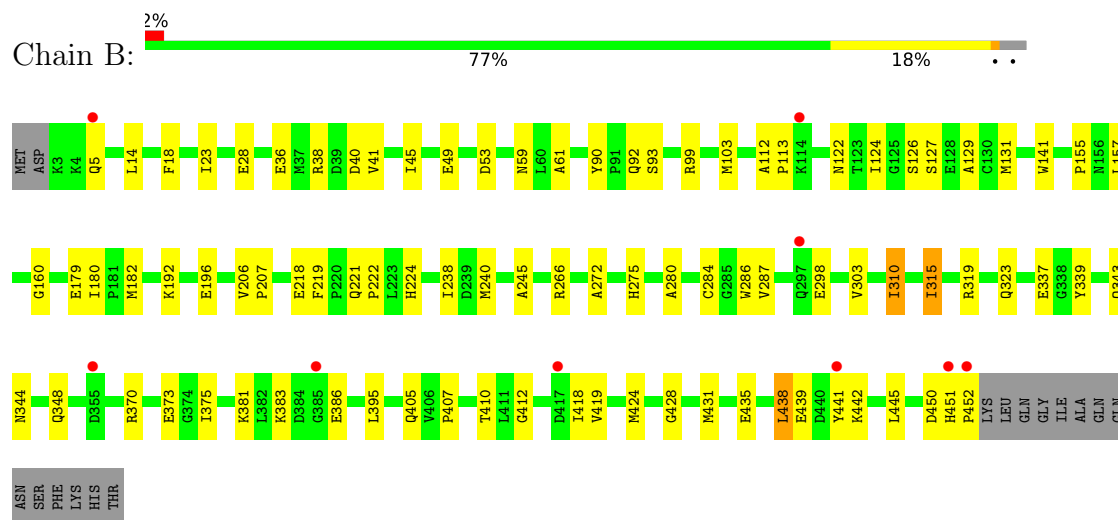
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

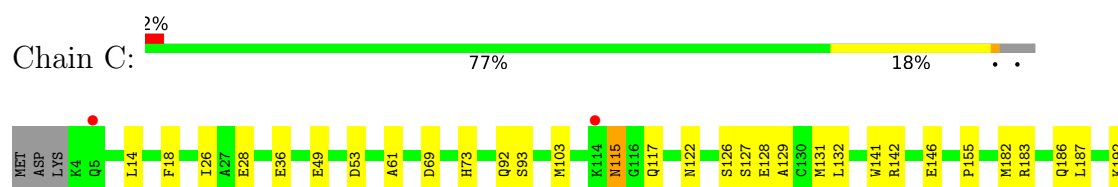
• Molecule 1: Glutamate decarboxylase beta

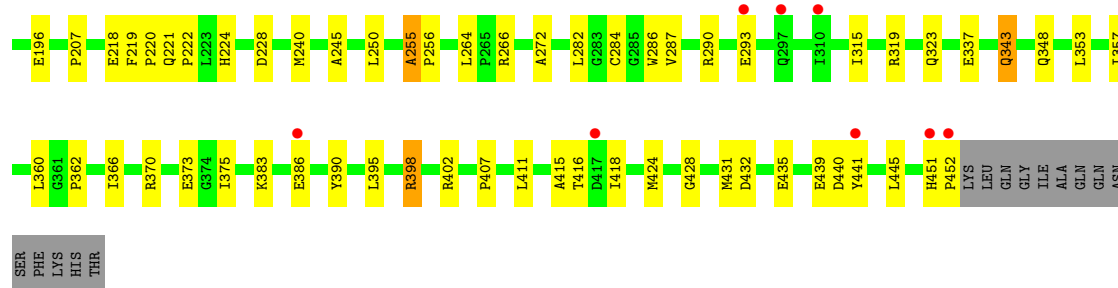


• Molecule 1: Glutamate decarboxylase beta

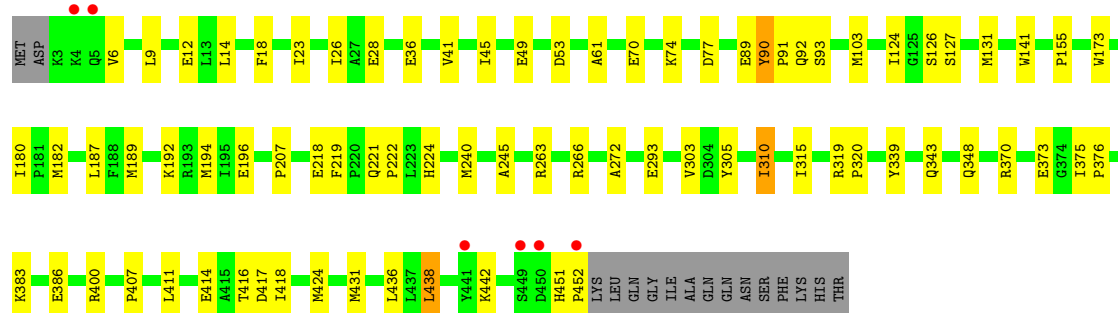
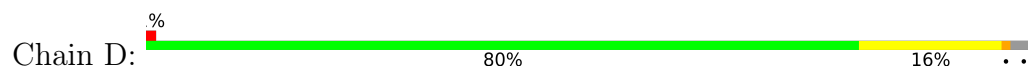


• Molecule 1: Glutamate decarboxylase beta

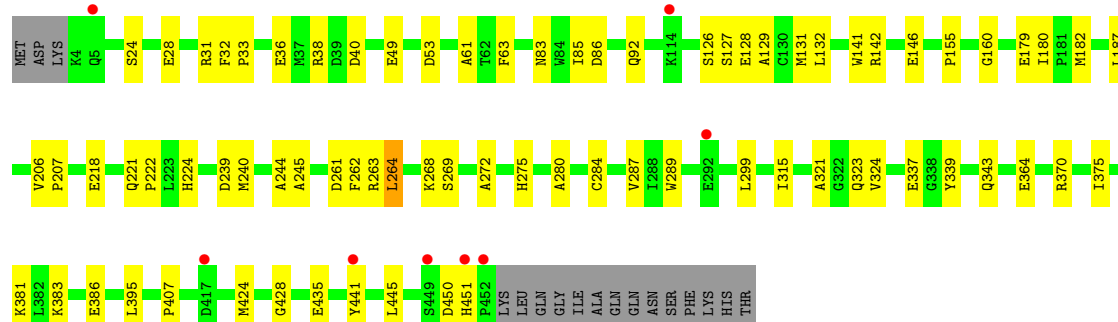
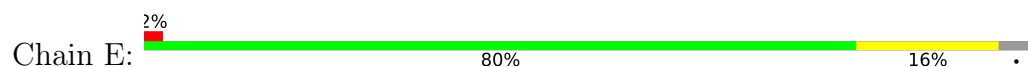




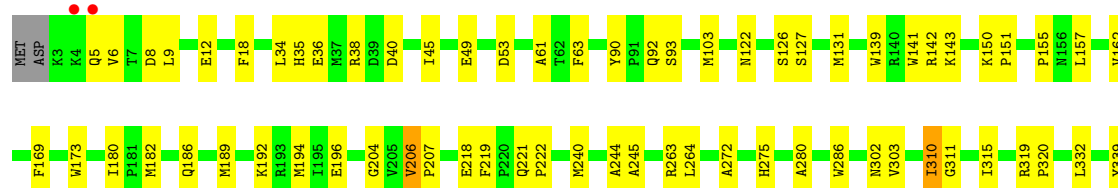
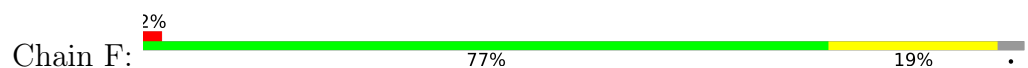
- Molecule 1: Glutamate decarboxylase beta



- Molecule 1: Glutamate decarboxylase beta



- Molecule 1: Glutamate decarboxylase beta





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	90.78Å 91.41Å 93.69Å 76.80° 77.10° 78.24°	Depositor
Resolution (Å)	24.62 – 2.00 24.62 – 2.01	Depositor EDS
% Data completeness (in resolution range)	95.8 (24.62-2.00) 95.2 (24.62-2.01)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.10 (at 2.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.182 , 0.213 0.181 , 0.211	Depositor DCC
R_{free} test set	5238 reflections (2.93%)	wwPDB-VP
Wilson B-factor (Å ²)	15.1	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.146 for k,l,h 0.146 for l,h,k 0.019 for -k,-h,-l 0.012 for -l,-k,-h 0.012 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23447	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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, ACY

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	0.36	0/3691	0.88	8/5002 (0.2%)
1	B	0.37	0/3692	0.87	8/5004 (0.2%)
1	C	0.35	0/3691	0.88	12/5004 (0.2%)
1	D	0.37	0/3691	0.88	7/5002 (0.1%)
1	E	0.36	0/3682	0.89	7/4991 (0.1%)
1	F	0.37	0/3691	0.88	11/5002 (0.2%)
All	All	0.37	0/22138	0.88	53/30005 (0.2%)

There are no bond length outliers.

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	218	GLU	N-CA-C	-7.95	95.78	108.73
1	A	218	GLU	N-CA-C	-7.71	96.15	108.73
1	F	218	GLU	N-CA-C	-7.67	96.23	108.73
1	D	218	GLU	N-CA-C	-7.59	96.36	108.73
1	E	218	GLU	N-CA-C	-7.57	96.39	108.73
1	D	90	TYR	CA-C-N	7.47	127.84	119.32
1	D	90	TYR	C-N-CA	7.47	127.84	119.32
1	C	218	GLU	N-CA-C	-7.36	96.74	108.73
1	F	93	SER	N-CA-C	-6.09	104.72	111.36
1	C	264	LEU	CA-C-N	6.08	125.56	119.24
1	C	264	LEU	C-N-CA	6.08	125.56	119.24
1	E	263	ARG	N-CA-C	-5.84	105.41	112.54
1	E	36	GLU	N-CA-C	-5.78	102.74	110.55
1	A	264	LEU	CA-C-N	5.74	125.21	119.24
1	A	264	LEU	C-N-CA	5.74	125.21	119.24
1	F	264	LEU	CA-C-N	5.69	125.15	119.24
1	F	264	LEU	C-N-CA	5.69	125.15	119.24
1	D	53	ASP	N-CA-C	-5.69	102.87	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	93	SER	N-CA-C	-5.68	105.09	111.28
1	D	93	SER	N-CA-C	-5.64	105.14	111.28
1	A	14	LEU	N-CA-C	5.63	120.22	113.23
1	C	36	GLU	N-CA-C	-5.59	103.00	110.55
1	A	93	SER	N-CA-C	-5.58	105.20	111.28
1	C	255	ALA	CA-C-N	5.58	125.94	119.47
1	C	255	ALA	C-N-CA	5.58	125.94	119.47
1	F	53	ASP	N-CA-C	-5.53	103.09	110.55
1	F	206	VAL	N-CA-C	5.51	113.61	108.15
1	E	206	VAL	N-CA-C	5.45	113.54	108.15
1	C	93	SER	N-CA-C	-5.44	105.35	111.28
1	C	14	LEU	N-CA-C	5.42	119.96	113.23
1	B	53	ASP	N-CA-C	-5.42	102.47	110.48
1	C	319	ARG	CA-C-N	5.41	125.33	119.76
1	C	319	ARG	C-N-CA	5.41	125.33	119.76
1	A	36	GLU	N-CA-C	-5.40	103.26	110.55
1	A	53	ASP	N-CA-C	-5.36	103.31	110.55
1	E	53	ASP	N-CA-C	-5.31	103.39	110.55
1	B	238	ILE	N-CA-C	5.27	115.55	108.17
1	B	90	TYR	CA-C-N	5.21	125.02	119.28
1	B	90	TYR	C-N-CA	5.21	125.02	119.28
1	C	366	ILE	N-CA-C	-5.21	107.33	111.81
1	F	90	TYR	CA-C-N	5.21	125.01	119.28
1	F	90	TYR	C-N-CA	5.21	125.01	119.28
1	F	162	VAL	N-CA-C	5.16	116.04	110.21
1	A	263	ARG	N-CA-C	-5.12	106.30	112.54
1	D	263	ARG	N-CA-C	-5.11	106.31	112.54
1	F	366	ILE	N-CA-C	-5.09	107.43	111.81
1	F	263	ARG	N-CA-C	-5.09	106.33	112.54
1	D	14	LEU	N-CA-C	5.09	119.54	113.23
1	B	315	ILE	N-CA-C	-5.08	106.19	111.58
1	E	264	LEU	CA-C-N	5.08	125.11	119.32
1	E	264	LEU	C-N-CA	5.08	125.11	119.32
1	B	14	LEU	N-CA-C	5.06	119.51	113.23
1	C	53	ASP	N-CA-C	-5.03	103.76	110.55

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	3584	0	3494	66	0
1	B	3585	0	3493	63	0
1	C	3579	0	3478	66	0
1	D	3584	0	3494	59	0
1	E	3575	0	3481	53	0
1	F	3584	0	3494	63	0
2	A	15	0	6	1	0
2	B	15	0	6	1	0
2	C	15	0	6	1	0
2	D	15	0	6	1	0
2	E	15	0	6	1	0
2	F	15	0	6	1	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
3	C	4	0	3	0	0
3	D	4	0	3	0	0
3	E	8	0	6	1	0
4	A	309	0	0	2	0
4	B	320	0	0	0	0
4	C	284	0	0	1	0
4	D	333	0	0	1	0
4	E	292	0	0	0	0
4	F	304	0	0	1	0
All	All	23447	0	20988	324	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 8.

All (324) 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:C:182:MET:HE2	1:C:187:LEU:HB3	1.50	0.94
1:F:139:TRP:HE1	1:F:143:LYS:HE3	1.33	0.93
1:B:23:ILE:HG13	1:B:45:ILE:HD11	1.61	0.82
1:D:23:ILE:HG13	1:D:45:ILE:HD11	1.60	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:139:TRP:NE1	1:F:143:LYS:HE3	1.94	0.81
1:C:131[B]:MET:SD	1:D:315:ILE:HG23	2.22	0.80
1:D:310:ILE:HD13	1:D:310:ILE:H	1.48	0.78
1:F:127:SER:O	1:F:131[B]:MET:HG2	1.85	0.76
1:E:131[B]:MET:SD	1:F:315:ILE:HG23	2.28	0.74
1:E:127:SER:O	1:E:131[A]:MET:HG2	1.86	0.74
1:A:182:MET:HE2	1:A:187:LEU:HB3	1.71	0.73
1:D:438:LEU:HD22	1:D:442:LYS:HE3	1.72	0.72
1:B:127:SER:O	1:B:131[B]:MET:HG2	1.90	0.71
1:E:395:LEU:HD21	1:E:441:TYR:CZ	2.25	0.71
1:A:131[B]:MET:SD	1:B:315:ILE:HG23	2.30	0.71
1:C:127:SER:O	1:C:131[A]:MET:HG2	1.91	0.71
1:B:370:ARG:HH11	1:B:370:ARG:HG2	1.56	0.70
1:F:348:GLN:HE21	1:F:431:MET:CE	2.03	0.70
1:F:339:TYR:O	1:F:343:GLN:HG2	1.91	0.70
1:B:221:GLN:HB3	1:B:222:PRO:HD3	1.74	0.69
1:E:339:TYR:O	1:E:343:GLN:HG2	1.92	0.69
1:D:339:TYR:O	1:D:343:GLN:HG2	1.93	0.69
1:D:127:SER:O	1:D:131[B]:MET:HG2	1.93	0.68
1:C:207:PRO:HG2	1:C:240:MET:HE2	1.75	0.68
1:C:395:LEU:HD21	1:C:441:TYR:CZ	2.27	0.68
1:F:395:LEU:HD21	1:F:441:TYR:CZ	2.29	0.68
1:C:192:LYS:O	1:C:196:GLU:HG3	1.94	0.68
1:E:221:GLN:HB3	1:E:222:PRO:HD3	1.76	0.68
1:A:395:LEU:HD21	1:A:441:TYR:CZ	2.28	0.68
1:B:339:TYR:O	1:B:343:GLN:HG2	1.94	0.68
1:F:131[B]:MET:HE3	1:F:173:TRP:HZ3	1.58	0.67
1:C:250:LEU:HD23	1:C:343:GLN:HG3	1.75	0.67
1:F:207:PRO:HG2	1:F:240:MET:HE2	1.78	0.66
1:A:127:SER:O	1:A:131[B]:MET:HG3	1.96	0.66
1:E:33:PRO:HD3	1:F:103:MET:HE2	1.76	0.66
1:C:142:ARG:O	1:C:146:GLU:HG3	1.96	0.66
1:B:438:LEU:HD22	1:B:442:LYS:HE3	1.78	0.65
1:A:383:LYS:HB2	1:A:386:GLU:HG3	1.77	0.65
1:A:127:SER:O	1:A:131[A]:MET:HG2	1.97	0.65
1:B:348[B]:GLN:HG2	1:B:431:MET:HE1	1.79	0.64
1:E:207:PRO:HG2	1:E:240:MET:HE2	1.78	0.64
1:A:375:ILE:HD12	1:A:424[A]:MET:HE1	1.78	0.63
1:D:89:GLU:C	1:D:91:PRO:HD3	2.24	0.63
1:A:339:TYR:O	1:A:343:GLN:HG2	1.97	0.63
1:A:26:ILE:HD13	1:F:435:GLU:HG3	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:207:PRO:HG2	1:D:240:MET:HE2	1.81	0.63
1:E:364:GLU:OE1	1:E:381:LYS:HE3	1.99	0.63
1:A:441:TYR:HE1	1:A:445:LEU:HD11	1.64	0.62
1:B:383:LYS:HB2	1:B:386:GLU:HG3	1.81	0.62
1:D:348:GLN:HE21	1:D:431:MET:HE2	1.65	0.61
1:E:375:ILE:HD12	1:E:424[A]:MET:HE1	1.80	0.61
1:E:32:PHE:HB2	4:F:741:HOH:O	2.00	0.61
1:B:344:ASN:O	1:B:348[A]:GLN:HG3	2.01	0.61
1:B:245:ALA:HA	1:B:272:ALA:HA	1.82	0.60
1:D:348:GLN:HE21	1:D:431:MET:CE	2.14	0.60
1:C:441:TYR:HE1	1:C:445:LEU:HD11	1.66	0.60
1:D:6:VAL:HG11	1:F:5:GLN:HG2	1.84	0.60
1:C:61:ALA:HB2	1:C:407:PRO:HD3	1.84	0.59
1:F:180:ILE:HD12	1:F:189:MET:HG3	1.83	0.59
1:C:183:ARG:HD2	1:C:186:GLN:OE1	2.02	0.59
1:A:126:SER:HB2	2:A:500:PLP:O4P	2.03	0.59
1:C:435:GLU:O	1:C:439:GLU:HG3	2.03	0.58
1:B:435:GLU:HG3	1:C:26:ILE:HD13	1.86	0.58
1:F:348:GLN:HE21	1:F:431:MET:HE3	1.68	0.58
1:D:305:TYR:HB2	1:D:310:ILE:HD12	1.85	0.58
1:E:28:GLU:HG2	1:E:31:ARG:O	2.02	0.58
1:A:207:PRO:HG2	1:A:240:MET:HE2	1.85	0.58
1:A:3:LYS:O	1:A:6:VAL:HG12	2.04	0.58
1:B:395:LEU:HD21	1:B:441:TYR:CZ	2.39	0.58
1:A:221:GLN:HB3	1:A:222:PRO:HD3	1.87	0.57
1:E:49:GLU:O	1:F:92:GLN:HG2	2.05	0.57
1:D:310:ILE:HD13	1:D:310:ILE:N	2.19	0.57
1:F:435:GLU:O	1:F:439:GLU:HG3	2.05	0.57
1:C:115:ASN:ND2	1:C:117:GLN:H	2.03	0.56
1:E:126:SER:HB2	2:E:502:PLP:O4P	2.05	0.56
1:D:303:VAL:HG23	1:D:310:ILE:HD11	1.86	0.56
1:C:221:GLN:HB3	1:C:222:PRO:HD3	1.87	0.56
1:F:139:TRP:CD1	1:F:143:LYS:HE3	2.41	0.56
1:D:124:ILE:HD11	1:D:319:ARG:HD2	1.87	0.56
1:A:131[B]:MET:HE2	1:A:173:TRP:HZ3	1.71	0.56
1:E:142:ARG:O	1:E:146:GLU:HG3	2.06	0.56
1:D:375:ILE:HD12	1:D:424[A]:MET:HE1	1.87	0.55
1:D:180:ILE:HD12	1:D:189:MET:HG3	1.88	0.55
1:A:70:GLU:HG2	1:A:74:LYS:HE3	1.88	0.55
1:F:182:MET:CE	1:F:412:GLY:H	2.19	0.55
1:A:61:ALA:HB2	1:A:407:PRO:HD3	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ILE:CD1	1:A:424[A]:MET:HE1	2.36	0.54
1:B:207:PRO:HG2	1:B:240:MET:HE2	1.89	0.54
1:C:395:LEU:HD21	1:C:441:TYR:CE2	2.42	0.54
1:E:315:ILE:HG21	1:F:315:ILE:HG21	1.90	0.54
1:D:418:ILE:HD12	1:D:418:ILE:N	2.22	0.54
1:A:441:TYR:CE1	1:A:445:LEU:HD11	2.43	0.54
1:E:92:GLN:HG2	1:F:49:GLU:O	2.07	0.54
1:A:370:ARG:HG2	1:A:370:ARG:HH11	1.73	0.54
1:F:410:THR:HG22	1:F:419:VAL:HG22	1.90	0.54
1:A:36:GLU:HG2	1:B:337:GLU:CD	2.33	0.54
1:F:8:ASP:O	1:F:12:GLU:HG3	2.08	0.53
1:B:370:ARG:HG2	1:B:370:ARG:NH1	2.23	0.53
1:D:126:SER:HB2	2:D:501:PLP:O4P	2.07	0.53
1:C:126:SER:HB2	2:C:501:PLP:O4P	2.08	0.53
1:D:305:TYR:CB	1:D:310:ILE:HD12	2.38	0.53
1:A:298:GLU:H	1:A:298:GLU:CD	2.16	0.53
1:A:315:ILE:HG21	1:B:315:ILE:HG21	1.90	0.53
1:A:395:LEU:HD21	1:A:441:TYR:CE2	2.44	0.53
1:C:245:ALA:HA	1:C:272:ALA:HA	1.89	0.53
1:E:239:ASP:CG	1:E:268:LYS:HZ1	2.16	0.53
1:F:180:ILE:HD11	1:F:194:MET:HA	1.91	0.53
1:F:126:SER:HB2	2:F:502:PLP:O4P	2.09	0.53
1:D:245:ALA:HA	1:D:272:ALA:HA	1.90	0.53
1:D:383:LYS:HB2	1:D:386:GLU:HG3	1.91	0.52
1:E:61:ALA:HB2	1:E:407:PRO:HD3	1.89	0.52
1:E:441:TYR:HE1	1:E:445:LEU:HD11	1.73	0.52
1:A:157:LEU:C	1:A:157:LEU:HD13	2.34	0.52
1:A:92:GLN:HG2	1:B:49:GLU:O	2.09	0.52
1:C:141:TRP:CZ3	1:C:155:PRO:HB3	2.45	0.52
1:B:348[B]:GLN:HG2	1:B:431:MET:CE	2.39	0.51
1:C:73:HIS:HD2	1:D:77:ASP:OD1	1.93	0.51
1:F:348:GLN:HE21	1:F:431:MET:HE2	1.75	0.51
1:D:293:GLU:HG3	4:D:7807:HOH:O	2.10	0.51
1:A:315:ILE:HG23	1:B:131[A]:MET:SD	2.51	0.51
1:E:129:ALA:HB1	1:E:287:VAL:HB	1.92	0.51
1:C:411:LEU:O	1:C:416:THR:HA	2.10	0.51
1:B:303:VAL:CG2	1:B:310:ILE:HG23	2.41	0.51
1:E:182:MET:HE2	1:E:187:LEU:HB3	1.93	0.50
1:F:375:ILE:HD12	1:F:424[B]:MET:HE1	1.93	0.50
1:A:128:GLU:O	1:A:132:LEU:HG	2.11	0.50
1:C:315:ILE:HD13	1:D:315:ILE:HG21	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:LYS:NZ	1:A:298:GLU:OE2	2.33	0.50
1:A:171:ARG:NH1	1:B:298:GLU:O	2.42	0.50
1:C:348:GLN:HG2	1:C:431:MET:HE1	1.92	0.50
1:C:402:ARG:HE	1:C:440:ASP:CG	2.19	0.50
1:B:160:GLY:O	1:B:179:GLU:HG3	2.12	0.50
1:B:129:ALA:HB1	1:B:287:VAL:HB	1.93	0.50
1:C:284:CYS:HB2	1:C:323:GLN:HB3	1.94	0.50
1:E:63:PHE:CE1	1:E:424[B]:MET:HE3	2.47	0.50
1:D:131[B]:MET:HE3	1:D:173:TRP:HZ3	1.77	0.50
1:A:245:ALA:HA	1:A:272:ALA:HA	1.94	0.49
1:C:250:LEU:CD2	1:C:343:GLN:HG3	2.42	0.49
1:D:221:GLN:HB3	1:D:222:PRO:HD3	1.95	0.49
1:B:126:SER:HB2	2:B:500:PLP:O4P	2.12	0.49
1:C:250:LEU:HD23	1:C:343:GLN:CG	2.40	0.49
1:C:375:ILE:HD12	1:C:424:MET:HE1	1.93	0.49
1:D:436:LEU:HD11	1:E:49:GLU:HB3	1.93	0.49
1:B:395:LEU:HD21	1:B:441:TYR:CE2	2.48	0.49
1:C:129:ALA:HB1	1:C:287:VAL:HB	1.94	0.49
1:D:180:ILE:HD11	1:D:194:MET:HA	1.94	0.49
1:E:28:GLU:HG3	1:F:103:MET:HE3	1.94	0.49
1:F:372:ASP:OD1	1:F:373:GLU:HG3	2.13	0.49
1:E:38:ARG:HB3	1:E:40:ASP:OD1	2.13	0.48
1:B:192:LYS:O	1:B:196:GLU:HG3	2.12	0.48
1:C:370:ARG:HD3	1:C:373:GLU:CD	2.38	0.48
1:C:441:TYR:CE1	1:C:445:LEU:HD11	2.48	0.48
1:F:63:PHE:CE1	1:F:424[A]:MET:HE3	2.48	0.48
1:D:182:MET:HE2	1:D:187:LEU:HB3	1.95	0.48
1:D:224:HIS:CD2	1:D:266:ARG:HB2	2.49	0.48
1:B:438:LEU:CD2	1:B:442:LYS:HE3	2.43	0.48
1:A:428:GLY:HA2	1:F:18:PHE:CE1	2.49	0.48
1:E:383:LYS:HB2	1:E:386:GLU:HG3	1.94	0.48
1:F:302:ASN:HD22	1:F:311:GLY:HA2	1.78	0.48
1:D:90:TYR:N	1:D:91:PRO:HD3	2.28	0.48
1:D:9:LEU:HG	1:F:9:LEU:HD21	1.94	0.47
1:E:321:ALA:O	1:E:324:VAL:HG12	2.14	0.47
1:F:245:ALA:HA	1:F:272:ALA:HA	1.95	0.47
1:A:17:ARG:HD3	1:A:44:GLN:HB2	1.96	0.47
1:A:219:PHE:O	1:A:222:PRO:HD2	2.14	0.47
1:A:141:TRP:CZ3	1:A:155:PRO:HB3	2.49	0.47
1:C:315:ILE:HG23	1:D:131[A]:MET:SD	2.54	0.47
1:F:141:TRP:CZ3	1:F:155:PRO:HB3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:207:PRO:CG	1:F:240:MET:HE2	2.42	0.47
1:C:255:ALA:N	1:C:256:PRO:HD3	2.30	0.47
1:C:115:ASN:C	1:C:115:ASN:HD22	2.22	0.47
1:F:186:GLN:O	1:F:186:GLN:HG3	2.15	0.47
1:A:36:GLU:HG2	1:B:337:GLU:OE1	2.14	0.47
1:A:446:LYS:HE3	1:A:450:ASP:OD2	2.15	0.47
1:A:129:ALA:HB1	1:A:287:VAL:HB	1.97	0.47
1:B:375:ILE:HD12	1:B:424:MET:HE1	1.95	0.47
1:C:370:ARG:HD3	1:C:373:GLU:OE1	2.15	0.47
1:E:141:TRP:CZ3	1:E:155:PRO:HB3	2.50	0.47
1:F:34:LEU:HD23	1:F:34:LEU:O	2.14	0.47
1:D:26:ILE:HD13	1:E:435:GLU:HG2	1.96	0.46
1:B:451:HIS:N	1:B:452:PRO:HD3	2.30	0.46
1:B:428:GLY:HA2	1:C:18:PHE:CE1	2.51	0.46
1:C:398:ARG:HH11	1:C:398:ARG:CG	2.29	0.46
1:E:284:CYS:HB2	1:E:323:GLN:HB3	1.98	0.46
1:C:128:GLU:O	1:C:132:LEU:HG	2.15	0.46
1:B:435:GLU:O	1:B:439:GLU:HG3	2.14	0.46
1:E:180:ILE:HD12	1:E:180:ILE:N	2.31	0.46
1:A:114:LYS:HG2	4:A:5816:HOH:O	2.15	0.45
1:C:337:GLU:CD	1:D:36:GLU:HG3	2.41	0.45
1:D:436:LEU:HD11	1:E:49:GLU:CB	2.46	0.45
1:E:395:LEU:HD21	1:E:441:TYR:CE2	2.51	0.45
1:B:141:TRP:CZ3	1:B:155:PRO:HB3	2.51	0.45
1:D:41:VAL:O	1:D:45:ILE:HG12	2.16	0.45
1:D:61:ALA:HB2	1:D:407:PRO:HD3	1.98	0.45
1:C:92:GLN:HG2	1:D:49:GLU:O	2.16	0.45
1:A:446:LYS:NZ	1:A:446:LYS:HA	2.31	0.45
1:C:228:ASP:HA	1:C:266:ARG:HH21	1.82	0.45
1:F:303:VAL:CG2	1:F:310:ILE:HG23	2.47	0.45
1:B:182:MET:CE	1:B:412:GLY:H	2.29	0.45
1:E:160:GLY:O	1:E:179:GLU:HG3	2.17	0.45
1:A:183:ARG:NH2	1:A:186:GLN:OE1	2.49	0.45
1:A:186:GLN:O	1:A:186:GLN:HG3	2.17	0.45
1:C:451:HIS:N	1:C:452:PRO:HD3	2.32	0.45
1:C:398:ARG:NH1	1:C:398:ARG:HG3	2.31	0.45
1:D:70:GLU:HG2	1:D:74:LYS:HE3	1.98	0.45
1:E:337:GLU:CD	1:F:36:GLU:HG3	2.42	0.45
1:E:450:ASP:C	1:E:451:HIS:ND1	2.75	0.45
1:A:337:GLU:CD	1:B:36:GLU:HG2	2.42	0.44
1:E:244:ALA:O	1:E:245:ALA:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:GLN:HG2	1:F:6:VAL:HG11	1.98	0.44
1:B:180:ILE:HD12	1:B:180:ILE:N	2.32	0.44
1:A:273:SER:CB	1:A:276:LYS:HG3	2.48	0.44
1:D:451:HIS:N	1:D:452:PRO:HD3	2.32	0.44
1:A:88:GLU:OE1	1:A:305:TYR:HA	2.18	0.44
1:B:219:PHE:O	1:B:222:PRO:HD2	2.18	0.44
1:C:343:GLN:OE1	1:C:343:GLN:HA	2.17	0.44
1:A:103:MET:HE2	1:B:28:GLU:HG3	1.99	0.44
1:A:446:LYS:HE3	1:A:450:ASP:CG	2.43	0.44
1:A:446:LYS:HA	1:A:446:LYS:HZ2	1.83	0.44
1:E:245:ALA:HA	1:E:272:ALA:HA	1.98	0.44
1:A:49:GLU:O	1:B:92:GLN:HG2	2.18	0.44
1:B:157:LEU:HD11	1:B:206:VAL:HG23	1.98	0.44
1:C:383:LYS:HB2	1:C:386:GLU:HG3	1.99	0.44
1:C:415:ALA:HB1	1:C:418:ILE:HD12	1.99	0.44
1:E:83:ASN:OD1	1:E:85:ILE:HG22	2.18	0.44
1:B:18:PHE:CE1	1:C:428:GLY:HA2	2.53	0.44
1:E:375:ILE:CD1	1:E:424[A]:MET:HE1	2.47	0.44
1:B:450:ASP:C	1:B:451:HIS:ND1	2.76	0.43
1:D:192:LYS:O	1:D:196:GLU:HG3	2.17	0.43
1:D:219:PHE:O	1:D:222:PRO:HD2	2.18	0.43
1:E:370:ARG:HH11	1:E:370:ARG:HG2	1.83	0.43
1:F:45:ILE:O	1:F:49:GLU:HG3	2.18	0.43
1:A:364:GLU:OE1	1:A:381:LYS:HE3	2.18	0.43
1:A:381:LYS:HD2	1:A:418:ILE:HG23	1.99	0.43
1:C:122:ASN:HB2	1:C:286:TRP:CZ3	2.53	0.43
1:C:362:PRO:HA	4:C:8790:HOH:O	2.17	0.43
1:F:61:ALA:HB2	1:F:407:PRO:HD3	1.98	0.43
1:B:38:ARG:HB3	1:B:40:ASP:OD1	2.19	0.43
1:B:61:ALA:HB2	1:B:407:PRO:HD3	1.99	0.43
1:C:28:GLU:HG3	1:D:103:MET:HE2	1.99	0.43
1:D:375:ILE:HB	1:D:376:PRO:CD	2.49	0.43
1:F:131[B]:MET:SD	1:F:169:PHE:HB2	2.58	0.43
1:A:273:SER:HB2	1:A:276:LYS:HG3	2.01	0.43
1:D:348:GLN:HE22	1:E:24:SER:HB2	1.83	0.43
1:D:141:TRP:CZ3	1:D:155:PRO:HB3	2.53	0.43
1:F:38:ARG:HB3	1:F:40:ASP:OD1	2.18	0.43
1:F:139:TRP:CE3	1:F:142:ARG:NH2	2.87	0.43
1:B:224:HIS:CD2	1:B:266:ARG:HB2	2.54	0.43
1:B:441:TYR:HE1	1:B:445:LEU:HD11	1.82	0.43
1:C:353:LEU:O	1:C:357:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:417:ASP:CG	1:D:418:ILE:HD12	2.44	0.43
1:B:284:CYS:HB2	1:B:323:GLN:HB3	1.99	0.43
1:D:411:LEU:O	1:D:416:THR:HA	2.19	0.43
1:A:157:LEU:HD22	1:A:204:GLY:O	2.19	0.43
1:C:219:PHE:O	1:C:222:PRO:HD2	2.19	0.43
1:E:343:GLN:OE1	1:E:343:GLN:HA	2.19	0.43
1:F:375:ILE:HB	1:F:376:PRO:CD	2.49	0.43
1:A:224:HIS:CD2	1:A:266:ARG:HB2	2.54	0.42
1:A:353:LEU:O	1:A:357:ILE:HG13	2.19	0.42
1:F:219:PHE:O	1:F:222:PRO:HD2	2.19	0.42
1:F:360:LEU:HB3	1:F:445:LEU:CD1	2.49	0.42
1:A:410:THR:HG22	1:A:419:VAL:HG22	2.00	0.42
1:B:59:ASN:HA	1:B:405:GLN:HB2	2.01	0.42
1:F:34:LEU:HD22	1:F:35:HIS:CE1	2.54	0.42
1:A:424[A]:MET:HG2	4:A:5632:HOH:O	2.19	0.42
1:A:180:ILE:N	1:A:180:ILE:HD12	2.34	0.42
1:C:219:PHE:HA	1:C:220:PRO:HD3	1.91	0.42
1:B:112:ALA:HA	1:B:113:PRO:HD3	1.86	0.42
1:A:63:PHE:CE1	1:A:424[B]:MET:HE3	2.54	0.42
1:B:23:ILE:CG1	1:B:45:ILE:HD11	2.39	0.42
1:B:410:THR:HG22	1:B:419:VAL:HG22	2.01	0.42
1:C:207:PRO:CG	1:C:240:MET:HE2	2.46	0.42
1:C:360:LEU:HB3	1:C:445:LEU:CD1	2.49	0.42
1:F:34:LEU:HD23	1:F:34:LEU:C	2.45	0.42
1:F:221:GLN:HB3	1:F:222:PRO:HD3	2.01	0.42
1:C:315:ILE:HG21	1:D:315:ILE:HG21	2.02	0.42
1:F:157:LEU:HD11	1:F:206:VAL:HG23	2.02	0.42
1:A:86:ASP:HB3	1:A:89:GLU:HB2	2.00	0.42
1:E:28:GLU:CG	1:E:31:ARG:O	2.67	0.42
1:F:157:LEU:HD12	1:F:204:GLY:O	2.20	0.42
1:B:124:ILE:HD11	1:B:319:ARG:HD2	2.02	0.41
1:C:390:TYR:CD1	1:C:390:TYR:C	2.97	0.41
1:E:315:ILE:HG23	1:F:131[A]:MET:SD	2.60	0.41
1:A:38:ARG:HB3	1:A:40:ASP:OD1	2.20	0.41
1:E:128:GLU:O	1:E:132:LEU:HG	2.21	0.41
1:E:441:TYR:CE1	1:E:445:LEU:HD11	2.54	0.41
1:C:103:MET:HE2	1:D:28:GLU:CG	2.50	0.41
1:E:275:HIS:HA	1:E:280:ALA:O	2.21	0.41
1:E:86:ASP:OD2	3:E:9518:ACY:OXT	2.39	0.41
1:B:122:ASN:HB2	1:B:286:TRP:CZ3	2.56	0.41
1:E:269:SER:HB3	1:E:289:TRP:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:ASN:HD22	1:A:344:ASN:HA	1.68	0.41
1:B:207:PRO:CG	1:B:240:MET:HE2	2.49	0.41
1:B:381:LYS:HD2	1:B:418:ILE:HG23	2.03	0.41
1:C:49:GLU:O	1:D:92:GLN:HG2	2.21	0.41
1:C:115:ASN:ND2	1:C:115:ASN:C	2.79	0.41
1:D:124:ILE:CD1	1:D:319:ARG:HD2	2.48	0.41
1:E:261:ASP:HB2	1:E:262:PHE:H	1.64	0.41
1:F:122:ASN:HB2	1:F:286:TRP:CZ3	2.56	0.41
1:F:150:LYS:HB3	1:F:151:PRO:HD2	2.03	0.41
1:F:244:ALA:O	1:F:245:ALA:C	2.64	0.41
1:C:28:GLU:CG	1:D:103:MET:HE2	2.51	0.40
1:F:319:ARG:HB2	1:F:320:PRO:HD2	2.03	0.40
1:A:49:GLU:HG2	1:F:436:LEU:HD11	2.03	0.40
1:A:284:CYS:HB2	1:A:323:GLN:HB3	2.04	0.40
1:B:41:VAL:O	1:B:45:ILE:HG12	2.21	0.40
1:B:99:ARG:O	1:B:103:MET:HG3	2.21	0.40
1:C:282:LEU:O	1:D:320:PRO:HG3	2.21	0.40
1:F:275:HIS:HA	1:F:280:ALA:O	2.21	0.40
1:A:432:ASP:O	1:A:436:LEU:HD23	2.21	0.40
1:A:432:ASP:OD2	1:F:49:GLU:OE2	2.40	0.40
1:B:441:TYR:CE1	1:B:445:LEU:HD11	2.56	0.40
1:C:186:GLN:HG3	1:C:186:GLN:O	2.21	0.40
1:C:224:HIS:CD2	1:C:266:ARG:HB2	2.56	0.40
1:D:370:ARG:HD3	1:D:373:GLU:CD	2.46	0.40
1:F:192:LYS:O	1:F:196:GLU:HG3	2.20	0.40
1:F:451:HIS:N	1:F:452:PRO:HD3	2.36	0.40
1:B:275:HIS:HA	1:B:280:ALA:O	2.22	0.40
1:B:370:ARG:HD3	1:B:373:GLU:CD	2.47	0.40
1:C:69:ASP:OD1	1:C:69:ASP:C	2.65	0.40
1:D:18:PHE:CE1	1:E:428:GLY:HA2	2.56	0.40
1:B:49:GLU:OE2	1:C:432:ASP:OD2	2.40	0.40
1:D:303:VAL:CG2	1:D:310:ILE:HD11	2.50	0.40
1:E:224:HIS:CD2	1:E:264:LEU:HB3	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	451/466 (97%)	440 (98%)	11 (2%)	0	100	100
1	B	451/466 (97%)	440 (98%)	11 (2%)	0	100	100
1	C	451/466 (97%)	439 (97%)	12 (3%)	0	100	100
1	D	451/466 (97%)	438 (97%)	13 (3%)	0	100	100
1	E	450/466 (97%)	439 (98%)	11 (2%)	0	100	100
1	F	451/466 (97%)	440 (98%)	11 (2%)	0	100	100
All	All	2705/2796 (97%)	2636 (97%)	69 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	379/390 (97%)	377 (100%)	2 (0%)	81	87
1	B	379/390 (97%)	377 (100%)	2 (0%)	81	87
1	C	379/390 (97%)	374 (99%)	5 (1%)	61	68
1	D	379/390 (97%)	374 (99%)	5 (1%)	61	68
1	E	378/390 (97%)	377 (100%)	1 (0%)	86	91
1	F	379/390 (97%)	375 (99%)	4 (1%)	65	73
All	All	2273/2340 (97%)	2254 (99%)	19 (1%)	73	80

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

Mol	Chain	Res	Type
1	A	298	GLU
1	A	446	LYS
1	B	310	ILE
1	B	438	LEU
1	C	115	ASN
1	C	290	ARG
1	C	293	GLU
1	C	343	GLN
1	C	398	ARG
1	D	12	GLU
1	D	310	ILE
1	D	400	ARG
1	D	414	GLU
1	D	438	LEU
1	E	299	LEU
1	F	310	ILE
1	F	332	LEU
1	F	414	GLU
1	F	417	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	35	HIS
1	A	81	ASN
1	A	316	ASN
1	A	344	ASN
1	B	35	HIS
1	B	81	ASN
1	B	109	HIS
1	B	302	ASN
1	B	316	ASN
1	B	344	ASN
1	C	5	GLN
1	C	73	HIS
1	C	81	ASN
1	C	115	ASN
1	C	297	GLN
1	C	344	ASN
1	C	348	GLN

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Mol	Chain	Res	Type
1	D	35	HIS
1	D	81	ASN
1	D	109	HIS
1	D	201	ASN
1	D	348	GLN
1	E	35	HIS
1	E	81	ASN
1	E	109	HIS
1	F	5	GLN
1	F	297	GLN
1	F	302	ASN
1	F	344	ASN
1	F	348	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 oligosaccharides 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	PLP	B	500	1	15,15,16	0.77	0	21,22,23	1.50	6 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PLP	C	501	1	15,15,16	0.85	0	21,22,23	1.49	6 (28%)
3	ACY	C	8518	-	3,3,3	1.23	0	3,3,3	1.65	1 (33%)
2	PLP	A	500	1	15,15,16	0.86	0	21,22,23	1.47	5 (23%)
3	ACY	E	9519	-	3,3,3	1.21	0	3,3,3	1.64	1 (33%)
2	PLP	E	502	1	15,15,16	0.83	0	21,22,23	1.50	6 (28%)
2	PLP	F	502	1	15,15,16	0.79	0	21,22,23	1.50	6 (28%)
3	ACY	E	9518	-	3,3,3	1.27	0	3,3,3	1.60	1 (33%)
3	ACY	D	7518	-	3,3,3	1.26	0	3,3,3	1.62	1 (33%)
3	ACY	A	5518	-	3,3,3	1.23	0	3,3,3	1.59	1 (33%)
3	ACY	B	6518	-	3,3,3	1.19	0	3,3,3	1.66	1 (33%)
2	PLP	D	501	1	15,15,16	0.79	0	21,22,23	1.50	6 (28%)

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	PLP	B	500	1	-	0/6/6/8	0/1/1/1
2	PLP	C	501	1	-	0/6/6/8	0/1/1/1
2	PLP	A	500	1	-	0/6/6/8	0/1/1/1
2	PLP	E	502	1	-	0/6/6/8	0/1/1/1
2	PLP	F	502	1	-	0/6/6/8	0/1/1/1
2	PLP	D	501	1	-	0/6/6/8	0/1/1/1

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	PLP	C4A-C4-C5	3.12	124.16	120.94
2	E	502	PLP	C4A-C4-C5	2.99	124.02	120.94
2	F	502	PLP	C4A-C4-C5	2.98	124.02	120.94
2	B	500	PLP	C4A-C4-C5	2.96	123.99	120.94
2	D	501	PLP	C4A-C4-C5	2.94	123.97	120.94
2	A	500	PLP	C4A-C4-C5	2.86	123.89	120.94
2	D	501	PLP	C3-C2-N1	-2.74	117.50	120.96
2	C	501	PLP	C3-C2-N1	-2.72	117.53	120.96
2	A	500	PLP	C3-C2-N1	-2.72	117.53	120.96
2	F	502	PLP	C3-C2-N1	-2.71	117.55	120.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	502	PLP	C3-C2-N1	-2.69	117.56	120.96
2	B	500	PLP	C3-C2-N1	-2.68	117.57	120.96
2	B	500	PLP	C6-N1-C2	2.62	123.95	119.20
2	D	501	PLP	C6-N1-C2	2.60	123.91	119.20
2	F	502	PLP	C6-N1-C2	2.58	123.88	119.20
2	C	501	PLP	C6-N1-C2	2.57	123.86	119.20
2	E	502	PLP	C6-N1-C2	2.57	123.86	119.20
2	A	500	PLP	C6-N1-C2	2.56	123.85	119.20
2	E	502	PLP	C2A-C2-C3	2.37	123.57	120.80
2	D	501	PLP	C2A-C2-C3	2.31	123.50	120.80
2	B	500	PLP	C2A-C2-C3	2.29	123.48	120.80
2	C	501	PLP	C2A-C2-C3	2.29	123.47	120.80
2	F	502	PLP	C2A-C2-C3	2.28	123.46	120.80
3	B	6518	ACY	O-C-CH3	-2.27	113.24	122.53
3	C	8518	ACY	O-C-CH3	-2.25	113.28	122.53
2	A	500	PLP	C2A-C2-C3	2.25	123.43	120.80
3	E	9519	ACY	O-C-CH3	-2.24	113.34	122.53
3	D	7518	ACY	O-C-CH3	-2.20	113.50	122.53
3	E	9518	ACY	O-C-CH3	-2.18	113.57	122.53
2	D	501	PLP	C4A-C4-C3	-2.18	116.88	120.52
2	B	500	PLP	C4A-C4-C3	-2.17	116.91	120.52
3	A	5518	ACY	O-C-CH3	-2.17	113.64	122.53
2	F	502	PLP	C4A-C4-C3	-2.15	116.94	120.52
2	C	501	PLP	C4A-C4-C3	-2.14	116.95	120.52
2	E	502	PLP	C4A-C4-C3	-2.14	116.96	120.52
2	C	501	PLP	O3P-P-O1P	2.07	118.89	110.83
2	D	501	PLP	O3P-P-O1P	2.06	118.84	110.83
2	E	502	PLP	O3P-P-O1P	2.04	118.78	110.83
2	A	500	PLP	C4A-C4-C3	-2.03	117.14	120.52
2	B	500	PLP	O3P-P-O1P	2.03	118.73	110.83
2	F	502	PLP	O3P-P-O1P	2.01	118.68	110.83

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
2	B	500	PLP	1	0
2	C	501	PLP	1	0
2	A	500	PLP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	502	PLP	1	0
2	F	502	PLP	1	0
3	E	9518	ACY	1	0
2	D	501	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	450/466 (96%)	-0.16	8 (1%) 67 67	7, 13, 25, 44	3 (0%)
1	B	450/466 (96%)	-0.20	9 (2%) 65 64	7, 13, 26, 44	3 (0%)
1	C	449/466 (96%)	-0.06	10 (2%) 62 61	8, 15, 26, 44	4 (0%)
1	D	450/466 (96%)	-0.18	6 (1%) 75 74	5, 13, 26, 42	3 (0%)
1	E	449/466 (96%)	-0.14	8 (1%) 67 67	7, 14, 26, 46	3 (0%)
1	F	450/466 (96%)	-0.18	9 (2%) 65 64	6, 13, 26, 44	3 (0%)
All	All	2698/2796 (96%)	-0.15	50 (1%) 66 66	5, 14, 26, 46	19 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	452	PRO	4.1
1	F	452	PRO	3.7
1	B	385	GLY	3.3
1	E	5	GLN	3.3
1	F	5	GLN	3.2
1	C	452	PRO	3.0
1	F	385	GLY	3.0
1	A	452	PRO	3.0
1	A	417	ASP	2.9
1	B	5	GLN	2.9
1	B	452	PRO	2.8
1	C	441	TYR	2.8
1	E	114	LYS	2.8
1	E	452	PRO	2.8
1	F	417	ASP	2.7
1	C	310	ILE	2.6
1	F	4	LYS	2.6
1	D	441	TYR	2.6
1	C	451	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	3	LYS	2.5
1	B	441	TYR	2.5
1	C	5	GLN	2.4
1	E	441	TYR	2.4
1	D	5	GLN	2.4
1	C	114	LYS	2.4
1	A	386	GLU	2.3
1	C	386	GLU	2.3
1	E	292	GLU	2.3
1	C	297	GLN	2.3
1	C	293	GLU	2.3
1	A	114	LYS	2.3
1	E	451	HIS	2.3
1	F	451	HIS	2.3
1	B	355	ASP	2.3
1	D	4	LYS	2.3
1	D	450	ASP	2.2
1	A	451	HIS	2.2
1	E	417	ASP	2.2
1	B	417	ASP	2.2
1	E	449	SER	2.1
1	C	417	ASP	2.1
1	F	389	GLY	2.1
1	B	114	LYS	2.1
1	F	441	TYR	2.1
1	B	451	HIS	2.1
1	A	139	TRP	2.1
1	B	297	GLN	2.1
1	F	448	LEU	2.0
1	D	449	SER	2.0
1	A	447	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	ACY	E	9518	4/4	0.94	0.11	19,20,20,20	0
3	ACY	E	9519	4/4	0.94	0.09	17,17,17,17	0
3	ACY	C	8518	4/4	0.95	0.07	17,18,18,18	0
3	ACY	A	5518	4/4	0.96	0.06	18,18,19,19	0
3	ACY	B	6518	4/4	0.96	0.09	14,15,15,15	0
2	PLP	A	500	15/16	0.96	0.07	11,12,13,13	0
3	ACY	D	7518	4/4	0.96	0.07	19,19,19,20	0
2	PLP	C	501	15/16	0.96	0.07	12,13,14,15	0
2	PLP	F	502	15/16	0.96	0.07	10,11,11,12	0
2	PLP	D	501	15/16	0.97	0.06	10,10,10,11	0
2	PLP	E	502	15/16	0.97	0.06	11,12,13,13	0
2	PLP	B	500	15/16	0.97	0.06	11,11,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.