



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 30, 2021 – 12:14 PM JST

PDB ID : 6L07
Title : Crystal structure of Escherichia coli phosphatidylserine decarboxylase (PE-bound form)
Authors : Watanabe, Y.; Watanabe, S.
Deposited on : 2019-09-26
Resolution : 3.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

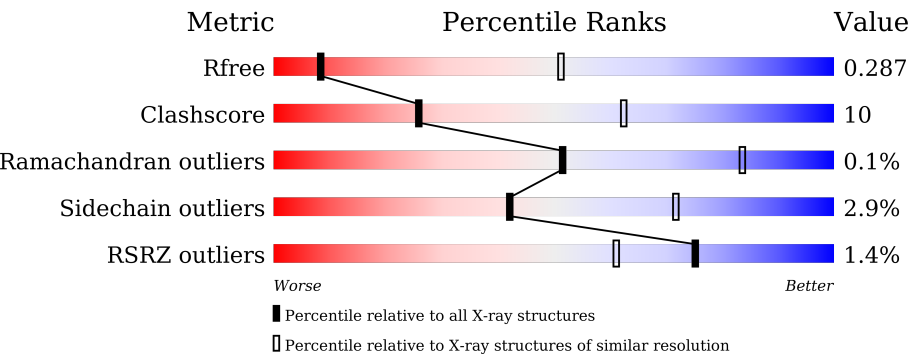
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 3.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1257 (3.70-3.50)
Clashscore	141614	1353 (3.70-3.50)
Ramachandran outliers	138981	1307 (3.70-3.50)
Sidechain outliers	138945	1307 (3.70-3.50)
RSRZ outliers	127900	1161 (3.70-3.50)

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	267	<div><div></div><div>69%26% . .</div></div>
1	B	267	<div><div></div><div>75%20% .</div></div>
1	C	267	<div><div></div><div>70%25% .</div></div>
1	D	267	<div><div></div><div>71%24% . .</div></div>
1	E	267	<div><div></div><div>%69%26% . .</div></div>
1	F	267	<div><div></div><div>69%26% . .</div></div>

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Mol	Chain	Length	Quality of chain
1	G	267	<div> <div>%</div> <div> <div></div> <div>64%</div> <div>31%</div> <div>• •</div> </div> </div>
1	H	267	<div> <div>6%</div> <div> <div></div> <div>69%</div> <div>26%</div> <div>•</div> </div> </div>
2	I	36	<div> <div></div> <div> <div></div> <div>67%</div> <div>31%</div> <div>•</div> </div> </div>
2	J	36	<div> <div></div> <div> <div></div> <div>78%</div> <div>19%</div> <div>•</div> </div> </div>
2	K	36	<div> <div></div> <div> <div></div> <div>69%</div> <div>28%</div> <div>•</div> </div> </div>
2	L	36	<div> <div></div> <div> <div></div> <div>78%</div> <div>19%</div> <div>•</div> </div> </div>
2	M	36	<div> <div></div> <div> <div></div> <div>64%</div> <div>31%</div> <div>6%</div> </div> </div>
3	N	36	<div> <div>3%</div> <div> <div></div> <div>61%</div> <div>36%</div> <div>•</div> </div> </div>
3	O	36	<div> <div>3%</div> <div> <div></div> <div>64%</div> <div>33%</div> <div>•</div> </div> </div>
3	P	36	<div> <div>14%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>•</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 18179 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 Phosphatidylserine decarboxylase beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			2025	1308	346	361	10			
1	B	255	Total	C	N	O	S	0	0	0
			2022	1307	346	359	10			
1	C	255	Total	C	N	O	S	0	0	0
			2005	1297	345	353	10			
1	D	255	Total	C	N	O	S	0	0	0
			2018	1304	344	360	10			
1	E	255	Total	C	N	O	S	0	0	0
			2013	1303	344	356	10			
1	F	255	Total	C	N	O	S	0	0	0
			2014	1302	346	357	9			
1	G	255	Total	C	N	O	S	0	0	0
			2022	1305	346	361	10			
1	H	255	Total	C	N	O	S	0	0	0
			1998	1289	343	356	10			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP A0A446DLT6
A	-12	GLY	-	expression tag	UNP A0A446DLT6
A	-11	SER	-	expression tag	UNP A0A446DLT6
A	-10	SER	-	expression tag	UNP A0A446DLT6
A	-9	HIS	-	expression tag	UNP A0A446DLT6
A	-8	HIS	-	expression tag	UNP A0A446DLT6
A	-7	HIS	-	expression tag	UNP A0A446DLT6
A	-6	HIS	-	expression tag	UNP A0A446DLT6
A	-5	HIS	-	expression tag	UNP A0A446DLT6
A	-4	HIS	-	expression tag	UNP A0A446DLT6
A	-3	SER	-	expression tag	UNP A0A446DLT6
A	-2	GLN	-	expression tag	UNP A0A446DLT6
A	-1	ASP	-	expression tag	UNP A0A446DLT6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	PRO	-	expression tag	UNP A0A446DLT6
B	-13	MET	-	expression tag	UNP A0A446DLT6
B	-12	GLY	-	expression tag	UNP A0A446DLT6
B	-11	SER	-	expression tag	UNP A0A446DLT6
B	-10	SER	-	expression tag	UNP A0A446DLT6
B	-9	HIS	-	expression tag	UNP A0A446DLT6
B	-8	HIS	-	expression tag	UNP A0A446DLT6
B	-7	HIS	-	expression tag	UNP A0A446DLT6
B	-6	HIS	-	expression tag	UNP A0A446DLT6
B	-5	HIS	-	expression tag	UNP A0A446DLT6
B	-4	HIS	-	expression tag	UNP A0A446DLT6
B	-3	SER	-	expression tag	UNP A0A446DLT6
B	-2	GLN	-	expression tag	UNP A0A446DLT6
B	-1	ASP	-	expression tag	UNP A0A446DLT6
B	0	PRO	-	expression tag	UNP A0A446DLT6
C	-13	MET	-	expression tag	UNP A0A446DLT6
C	-12	GLY	-	expression tag	UNP A0A446DLT6
C	-11	SER	-	expression tag	UNP A0A446DLT6
C	-10	SER	-	expression tag	UNP A0A446DLT6
C	-9	HIS	-	expression tag	UNP A0A446DLT6
C	-8	HIS	-	expression tag	UNP A0A446DLT6
C	-7	HIS	-	expression tag	UNP A0A446DLT6
C	-6	HIS	-	expression tag	UNP A0A446DLT6
C	-5	HIS	-	expression tag	UNP A0A446DLT6
C	-4	HIS	-	expression tag	UNP A0A446DLT6
C	-3	SER	-	expression tag	UNP A0A446DLT6
C	-2	GLN	-	expression tag	UNP A0A446DLT6
C	-1	ASP	-	expression tag	UNP A0A446DLT6
C	0	PRO	-	expression tag	UNP A0A446DLT6
D	-13	MET	-	expression tag	UNP A0A446DLT6
D	-12	GLY	-	expression tag	UNP A0A446DLT6
D	-11	SER	-	expression tag	UNP A0A446DLT6
D	-10	SER	-	expression tag	UNP A0A446DLT6
D	-9	HIS	-	expression tag	UNP A0A446DLT6
D	-8	HIS	-	expression tag	UNP A0A446DLT6
D	-7	HIS	-	expression tag	UNP A0A446DLT6
D	-6	HIS	-	expression tag	UNP A0A446DLT6
D	-5	HIS	-	expression tag	UNP A0A446DLT6
D	-4	HIS	-	expression tag	UNP A0A446DLT6
D	-3	SER	-	expression tag	UNP A0A446DLT6
D	-2	GLN	-	expression tag	UNP A0A446DLT6
D	-1	ASP	-	expression tag	UNP A0A446DLT6

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	PRO	-	expression tag	UNP A0A446DLT6
E	-13	MET	-	expression tag	UNP A0A446DLT6
E	-12	GLY	-	expression tag	UNP A0A446DLT6
E	-11	SER	-	expression tag	UNP A0A446DLT6
E	-10	SER	-	expression tag	UNP A0A446DLT6
E	-9	HIS	-	expression tag	UNP A0A446DLT6
E	-8	HIS	-	expression tag	UNP A0A446DLT6
E	-7	HIS	-	expression tag	UNP A0A446DLT6
E	-6	HIS	-	expression tag	UNP A0A446DLT6
E	-5	HIS	-	expression tag	UNP A0A446DLT6
E	-4	HIS	-	expression tag	UNP A0A446DLT6
E	-3	SER	-	expression tag	UNP A0A446DLT6
E	-2	GLN	-	expression tag	UNP A0A446DLT6
E	-1	ASP	-	expression tag	UNP A0A446DLT6
E	0	PRO	-	expression tag	UNP A0A446DLT6
F	-13	MET	-	expression tag	UNP A0A446DLT6
F	-12	GLY	-	expression tag	UNP A0A446DLT6
F	-11	SER	-	expression tag	UNP A0A446DLT6
F	-10	SER	-	expression tag	UNP A0A446DLT6
F	-9	HIS	-	expression tag	UNP A0A446DLT6
F	-8	HIS	-	expression tag	UNP A0A446DLT6
F	-7	HIS	-	expression tag	UNP A0A446DLT6
F	-6	HIS	-	expression tag	UNP A0A446DLT6
F	-5	HIS	-	expression tag	UNP A0A446DLT6
F	-4	HIS	-	expression tag	UNP A0A446DLT6
F	-3	SER	-	expression tag	UNP A0A446DLT6
F	-2	GLN	-	expression tag	UNP A0A446DLT6
F	-1	ASP	-	expression tag	UNP A0A446DLT6
F	0	PRO	-	expression tag	UNP A0A446DLT6
G	-13	MET	-	expression tag	UNP A0A446DLT6
G	-12	GLY	-	expression tag	UNP A0A446DLT6
G	-11	SER	-	expression tag	UNP A0A446DLT6
G	-10	SER	-	expression tag	UNP A0A446DLT6
G	-9	HIS	-	expression tag	UNP A0A446DLT6
G	-8	HIS	-	expression tag	UNP A0A446DLT6
G	-7	HIS	-	expression tag	UNP A0A446DLT6
G	-6	HIS	-	expression tag	UNP A0A446DLT6
G	-5	HIS	-	expression tag	UNP A0A446DLT6
G	-4	HIS	-	expression tag	UNP A0A446DLT6
G	-3	SER	-	expression tag	UNP A0A446DLT6
G	-2	GLN	-	expression tag	UNP A0A446DLT6
G	-1	ASP	-	expression tag	UNP A0A446DLT6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	PRO	-	expression tag	UNP A0A446DLT6
H	-13	MET	-	expression tag	UNP A0A446DLT6
H	-12	GLY	-	expression tag	UNP A0A446DLT6
H	-11	SER	-	expression tag	UNP A0A446DLT6
H	-10	SER	-	expression tag	UNP A0A446DLT6
H	-9	HIS	-	expression tag	UNP A0A446DLT6
H	-8	HIS	-	expression tag	UNP A0A446DLT6
H	-7	HIS	-	expression tag	UNP A0A446DLT6
H	-6	HIS	-	expression tag	UNP A0A446DLT6
H	-5	HIS	-	expression tag	UNP A0A446DLT6
H	-4	HIS	-	expression tag	UNP A0A446DLT6
H	-3	SER	-	expression tag	UNP A0A446DLT6
H	-2	GLN	-	expression tag	UNP A0A446DLT6
H	-1	ASP	-	expression tag	UNP A0A446DLT6
H	0	PRO	-	expression tag	UNP A0A446DLT6

- Molecule 2 is a protein called Phosphatidylserine decarboxylase alpha chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	35	Total	C	N	O	0	0	0
			244	159	39	46			
2	J	35	Total	C	N	O	0	0	0
			244	159	39	46			
2	K	35	Total	C	N	O	0	0	0
			244	159	39	46			
2	L	35	Total	C	N	O	0	0	0
			244	159	39	46			
2	M	34	Total	C	N	O	0	0	0
			239	156	38	45			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	254	PPI	SER	modified residue	UNP A0A446DLT6
J	254	PPI	SER	modified residue	UNP A0A446DLT6
K	254	PPI	SER	modified residue	UNP A0A446DLT6
L	254	PPI	SER	modified residue	UNP A0A446DLT6
M	254	PPI	SER	modified residue	UNP A0A446DLT6

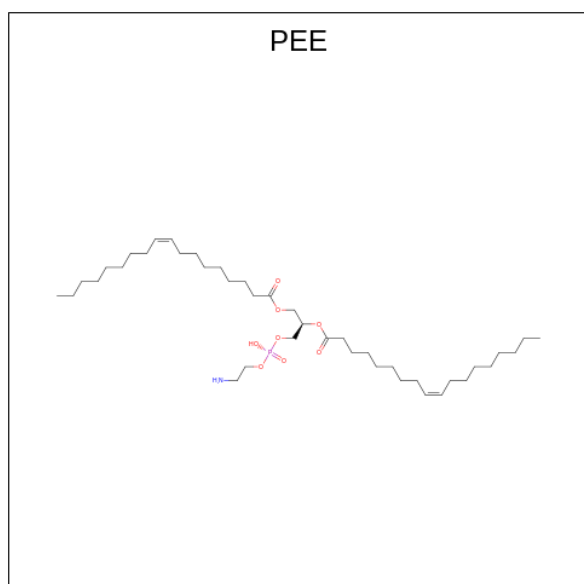
- Molecule 3 is a protein called Phosphatidylserine decarboxylase alpha chain.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	N	35	Total	C	N	O	0	0	0
			245	159	39	47			
3	O	35	Total	C	N	O	0	0	0
			245	159	39	47			
3	P	35	Total	C	N	O	0	0	0
			242	156	39	47			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	254	PYR	SER	modified residue	UNP A0A446DLT6
O	254	PYR	SER	modified residue	UNP A0A446DLT6
P	254	PYR	SER	modified residue	UNP A0A446DLT6

- Molecule 4 is 1,2-dioleoyl-sn-glycero-3-phosphoethanolamine (three-letter code: PEE) (formula: $C_{41}H_{78}NO_8P$).

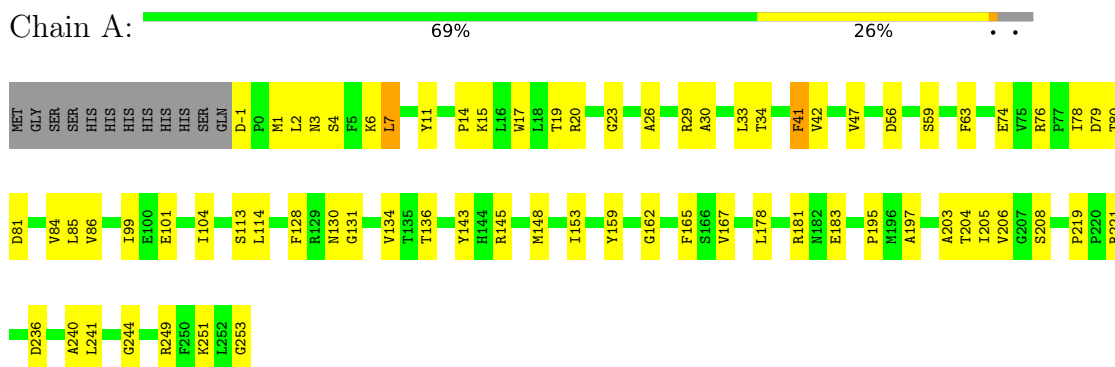


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	I	1	Total	C	N	O	P	0	0
			23	13	1	8	1		
4	J	1	Total	C	N	O	P	0	0
			23	13	1	8	1		
4	K	1	Total	C	N	O	P	0	0
			23	13	1	8	1		
4	L	1	Total	C	N	O	P	0	0
			23	13	1	8	1		
4	M	1	Total	C	N	O	P	0	0
			23	13	1	8	1		

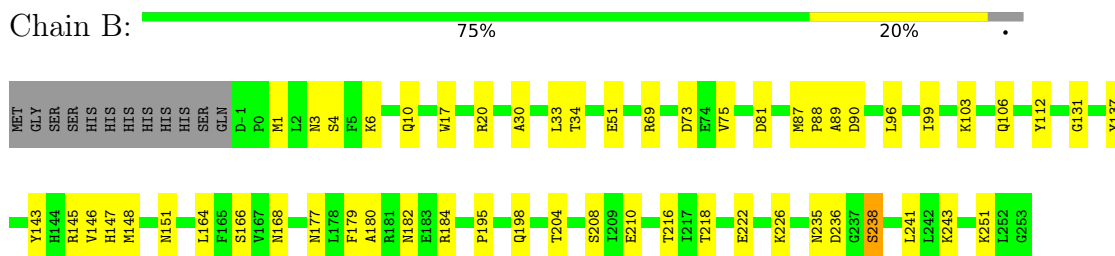
3 Residue-property plots

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.

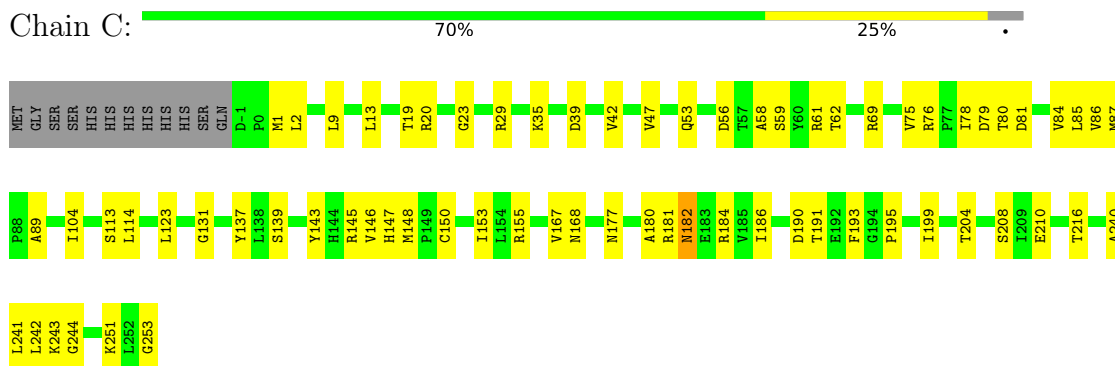
- Molecule 1: Phosphatidylserine decarboxylase beta chain



- Molecule 1: Phosphatidylserine decarboxylase beta chain

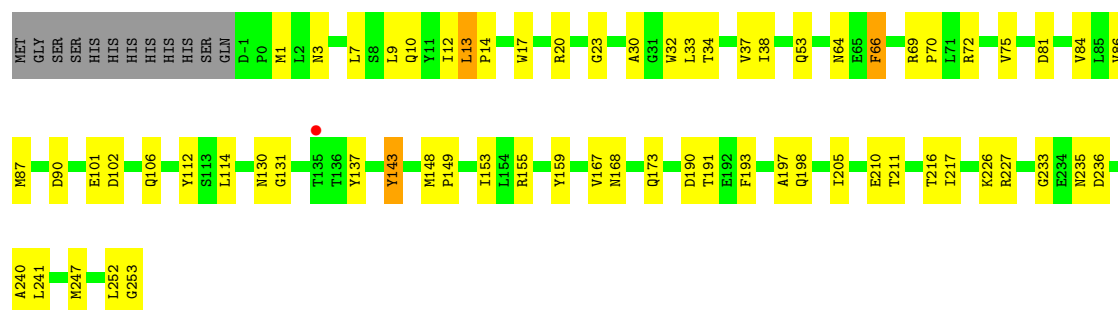


- Molecule 1: Phosphatidylserine decarboxylase beta chain



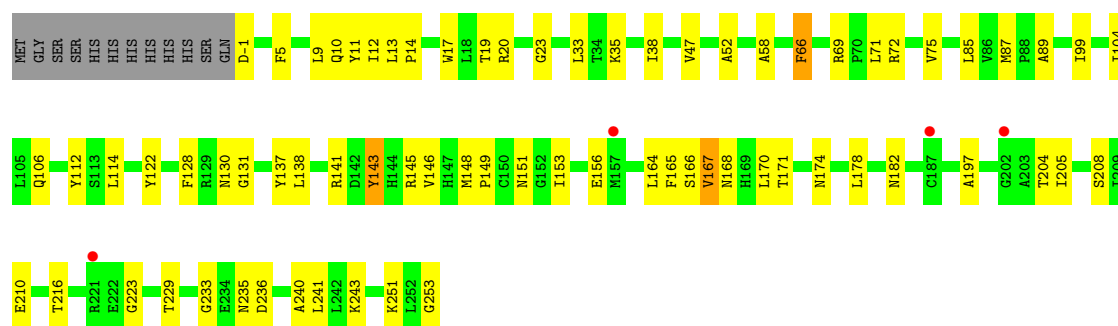
- Molecule 1: Phosphatidylserine decarboxylase beta chain

Chain D:  71% 24% . .



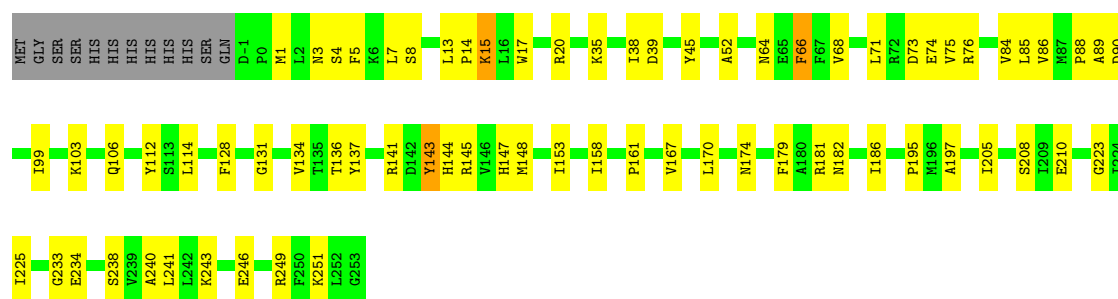
• Molecule 1: Phosphatidylserine decarboxylase beta chain

Chain E:  69% 26% . .



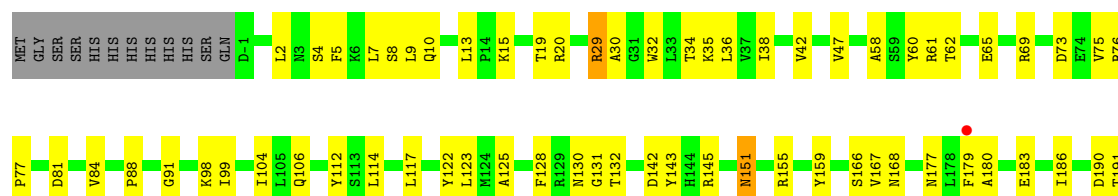
• Molecule 1: Phosphatidylserine decarboxylase beta chain

Chain F:  69% 26% . .



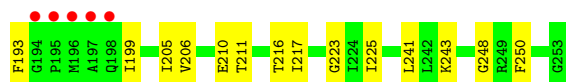
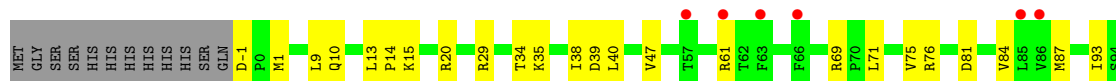
• Molecule 1: Phosphatidylserine decarboxylase beta chain

Chain G:  64% 31% . .

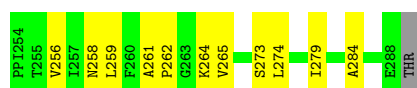




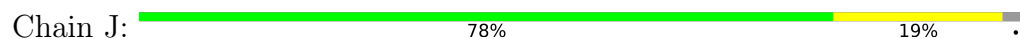
- Molecule 1: Phosphatidylserine decarboxylase beta chain



- Molecule 2: Phosphatidylserine decarboxylase alpha chain



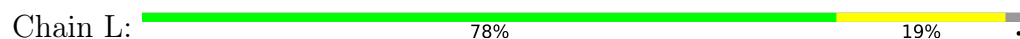
- Molecule 2: Phosphatidylserine decarboxylase alpha chain



- Molecule 2: Phosphatidylserine decarboxylase alpha chain

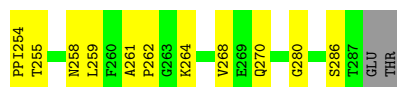


- Molecule 2: Phosphatidylserine decarboxylase alpha chain



- Molecule 2: Phosphatidylserine decarboxylase alpha chain

Chain M:  64% 31% 6%



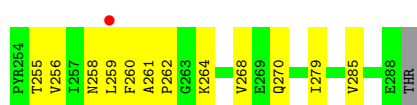
- Molecule 3: Phosphatidylserine decarboxylase alpha chain

Chain N:  3% 61% 36% .




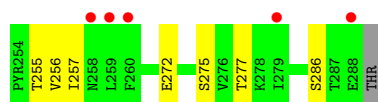
- Molecule 3: Phosphatidylserine decarboxylase alpha chain

Chain O:  3% 64% 33% .



- Molecule 3: Phosphatidylserine decarboxylase alpha chain

Chain P:  14% 78% 19% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	218.19Å 218.19Å 143.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.24 – 3.60 49.24 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.24-3.60) 99.9 (49.24-3.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 3.57Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.241 , 0.289 0.239 , 0.287	Depositor DCC
R_{free} test set	2328 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	160.1	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 98.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.044 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18179	wwPDB-VP
Average B, all atoms (Å ²)	154.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: PPI, PEE, PYR

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	0.27	0/2076	0.49	0/2819
1	B	0.25	0/2073	0.45	0/2815
1	C	0.25	0/2056	0.44	0/2795
1	D	0.24	0/2069	0.46	0/2811
1	E	0.25	0/2064	0.47	0/2805
1	F	0.24	0/2065	0.47	0/2806
1	G	0.25	0/2073	0.46	0/2815
1	H	0.24	0/2048	0.45	0/2785
2	I	0.24	0/242	0.50	0/331
2	J	0.24	0/242	0.49	0/331
2	K	0.25	0/242	0.46	0/331
2	L	0.24	0/242	0.50	0/331
2	M	0.26	0/237	0.51	0/324
3	N	0.24	0/242	0.49	0/331
3	O	0.23	0/242	0.45	0/331
3	P	0.23	0/239	0.46	0/327
All	All	0.25	0/18452	0.46	0/25088

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2025	0	2036	54	0
1	B	2022	0	2034	32	0
1	C	2005	0	2000	44	0
1	D	2018	0	2021	43	0
1	E	2013	0	2013	50	0
1	F	2014	0	2021	47	0
1	G	2022	0	2027	59	0
1	H	1998	0	1988	47	0
2	I	244	0	258	9	0
2	J	244	0	258	10	0
2	K	244	0	258	9	0
2	L	244	0	258	6	0
2	M	239	0	256	11	0
3	N	245	0	257	11	0
3	O	245	0	257	15	0
3	P	242	0	248	9	0
4	I	23	0	18	7	0
4	J	23	0	19	1	0
4	K	23	0	18	3	0
4	L	23	0	18	5	0
4	M	23	0	18	3	0
All	All	18179	0	18281	382	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 10.

All (382) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:ILE:HD13	1:A:104:ILE:HG12	1.54	0.88
1:A:131:GLY:HA2	2:I:262:PRO:HA	1.54	0.86
1:A:205:ILE:HD11	4:I:301:PEE:H49	1.62	0.80
1:A:30:ALA:HB3	1:A:34:THR:HG23	1.66	0.77
1:H:135:THR:HA	3:P:257:ILE:HG22	1.65	0.76
1:D:210:GLU:HB3	1:D:216:THR:HA	1.67	0.75
1:C:181:ARG:HD3	1:G:223:GLY:HA3	1.69	0.74
1:E:168:ASN:HD21	1:E:170:LEU:HB2	1.53	0.73
1:E:10:GLN:NE2	1:E:166:SER:OG	2.21	0.73
1:A:41:PHE:HD2	1:A:63:PHE:HZ	1.34	0.72
1:C:195:PRO:HD2	2:K:261:ALA:HB2	1.71	0.72
1:A:19:THR:HG23	1:A:205:ILE:HA	1.70	0.72
1:A:30:ALA:HB1	1:A:33:LEU:HB3	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:PRO:HD2	2:I:261:ALA:HB2	1.73	0.71
1:E:210:GLU:HB3	1:E:216:THR:HA	1.73	0.70
1:A:208:SER:HB3	1:A:251:LYS:HB3	1.72	0.70
1:E:89:ALA:HA	2:L:279:ILE:HG12	1.71	0.70
1:F:234:GLU:HG2	1:G:61:ARG:HH12	1.55	0.69
1:G:155:ARG:NH1	1:G:190:ASP:OD2	2.25	0.69
1:E:131:GLY:HA2	2:L:262:PRO:HA	1.74	0.69
1:E:19:THR:HG21	1:E:164:LEU:HD13	1.74	0.69
1:E:170:LEU:O	1:E:174:ASN:ND2	2.24	0.69
1:H:29:ARG:HH21	1:H:61:ARG:HD3	1.56	0.68
1:F:181:ARG:HD3	1:H:223:GLY:HA3	1.74	0.68
1:G:35:LYS:HE2	1:G:58:ALA:HB2	1.73	0.68
1:C:29:ARG:HH21	1:C:61:ARG:HD3	1.59	0.68
1:G:205:ILE:HG22	1:G:252:LEU:HG	1.75	0.68
1:H:98:LYS:HA	1:H:132:THR:HA	1.77	0.67
1:E:165:PHE:H	1:E:182:ASN:HD21	1.43	0.67
1:F:131:GLY:HA2	3:O:262:PRO:HA	1.78	0.66
1:C:186:ILE:HG23	1:C:199:ILE:HG12	1.76	0.66
1:E:166:SER:HB2	1:E:171:THR:HG21	1.78	0.66
1:D:233:GLY:O	1:D:235:ASN:ND2	2.29	0.66
1:G:10:GLN:NE2	1:G:166:SER:OG	2.27	0.66
1:H:126:ASP:HA	1:H:129:ARG:HD3	1.77	0.66
1:D:155:ARG:NH1	1:D:190:ASP:OD2	2.27	0.65
1:G:191:THR:HG23	1:G:193:PHE:H	1.61	0.65
1:D:14:PRO:HB2	1:D:17:TRP:HB2	1.79	0.65
1:C:199:ILE:HB	2:K:257:ILE:HG22	1.78	0.64
1:C:35:LYS:HE2	1:C:58:ALA:HB2	1.78	0.64
1:F:197:ALA:HB3	3:O:259:LEU:HB2	1.78	0.64
1:G:197:ALA:HB3	2:M:259:LEU:HB2	1.79	0.64
1:B:131:GLY:HA2	2:J:262:PRO:HA	1.80	0.64
1:C:210:GLU:HB3	1:C:216:THR:HA	1.80	0.64
4:I:301:PEE:O1P	4:I:301:PEE:H7	1.98	0.64
1:C:1:MET:HG3	1:C:2:LEU:HD12	1.80	0.63
1:B:30:ALA:HB3	1:B:34:THR:HG23	1.81	0.63
1:D:86:VAL:HG22	3:N:279:ILE:O	1.98	0.63
1:C:155:ARG:NH1	1:C:190:ASP:OD2	2.27	0.62
1:A:181:ARG:HD3	1:E:223:GLY:HA3	1.81	0.62
1:A:101:GLU:OE2	1:D:227:ARG:NH1	2.32	0.62
1:G:42:VAL:HG13	1:G:47:VAL:HB	1.79	0.62
1:G:30:ALA:O	1:G:34:THR:OG1	2.16	0.62
1:D:137:TYR:HE1	1:D:167:VAL:HG13	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:191:THR:HG23	1:D:193:PHE:H	1.65	0.61
1:C:123:LEU:HD11	1:G:123:LEU:HD11	1.82	0.61
1:C:131:GLY:HA2	2:K:262:PRO:HA	1.82	0.61
1:G:29:ARG:HH21	1:G:61:ARG:HH11	1.49	0.61
1:C:177:ASN:HB3	1:C:180:ALA:HB3	1.83	0.60
1:E:23:GLY:HA3	1:E:253:GLY:HA2	1.83	0.60
1:E:197:ALA:HB3	2:L:259:LEU:HB2	1.82	0.60
1:H:155:ARG:NH1	1:H:190:ASP:OD2	2.35	0.60
1:F:223:GLY:HA3	1:H:181:ARG:HD3	1.82	0.60
1:C:89:ALA:HB2	1:C:146:VAL:HG13	1.83	0.60
1:D:87:MET:HA	3:N:283:LEU:HD22	1.83	0.60
1:A:14:PRO:HB2	1:A:17:TRP:HD1	1.66	0.60
1:D:30:ALA:HB3	1:D:34:THR:HG23	1.84	0.60
4:I:301:PEE:H7	4:I:301:PEE:P	2.41	0.59
1:B:198:GLN:OE1	2:J:258:ASN:ND2	2.35	0.59
1:G:9:LEU:HG	1:G:13:LEU:HD23	1.84	0.59
1:C:147:HIS:ND1	1:C:244:GLY:O	2.34	0.59
1:D:131:GLY:HA2	3:N:262:PRO:HA	1.83	0.59
1:C:19:THR:OG1	1:C:204:THR:O	2.19	0.59
1:C:182:ASN:O	1:C:184:ARG:NH1	2.36	0.58
1:C:204:THR:HG22	4:K:301:PEE:H1	1.84	0.58
1:F:90:ASP:OD1	3:O:279:ILE:HG13	2.04	0.58
1:C:208:SER:HB3	1:C:251:LYS:HB3	1.85	0.58
1:A:1:MET:HA	1:A:4:SER:HB3	1.86	0.58
2:M:261:ALA:HB3	2:M:264:LYS:HG2	1.85	0.58
1:C:204:THR:HG21	4:K:301:PEE:H49	1.86	0.58
1:E:35:LYS:HE2	1:E:58:ALA:HB2	1.86	0.58
1:A:205:ILE:CD1	4:I:301:PEE:H49	2.33	0.58
1:G:98:LYS:HA	1:G:132:THR:HA	1.86	0.58
1:A:56:ASP:O	1:A:59:SER:OG	2.20	0.58
1:F:153:ILE:HG13	1:F:240:ALA:HB2	1.86	0.57
1:H:71:LEU:HD22	1:H:75:VAL:HG11	1.86	0.57
1:G:222:GLU:OE1	1:G:226:LYS:NZ	2.28	0.57
1:E:14:PRO:HB2	1:E:17:TRP:HD1	1.69	0.57
1:D:9:LEU:O	1:D:13:LEU:HG	2.05	0.57
1:H:191:THR:HG23	1:H:193:PHE:H	1.70	0.56
1:G:81:ASP:HB3	1:G:84:VAL:HG22	1.86	0.56
1:G:177:ASN:HB3	1:G:180:ALA:HB3	1.85	0.56
1:C:56:ASP:O	1:C:59:SER:OG	2.23	0.56
1:A:3:ASN:O	1:A:6:LYS:N	2.37	0.56
1:F:137:TYR:HE1	1:F:167:VAL:HG13	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:LYS:NZ	1:C:39:ASP:OD2	2.34	0.56
1:D:205:ILE:HG22	1:D:252:LEU:HG	1.88	0.56
1:F:89:ALA:HA	3:O:279:ILE:HG12	1.88	0.56
1:D:87:MET:HG2	3:N:281:GLN:O	2.06	0.55
1:G:208:SER:HB3	1:G:251:LYS:HB3	1.88	0.55
1:A:30:ALA:O	1:A:34:THR:OG1	2.19	0.55
1:E:10:GLN:OE1	1:E:166:SER:N	2.34	0.55
1:G:30:ALA:HB3	1:G:34:THR:HG23	1.87	0.55
1:B:182:ASN:O	1:B:184:ARG:NH1	2.40	0.55
1:C:104:ILE:HD11	1:C:114:LEU:HD13	1.88	0.55
1:C:191:THR:HG23	1:C:193:PHE:H	1.73	0.54
1:G:104:ILE:HD11	1:G:114:LEU:HD13	1.89	0.54
1:F:99:ILE:HG13	1:F:103:LYS:O	2.08	0.54
1:H:75:VAL:HG13	1:H:76:ARG:HG2	1.88	0.54
1:A:23:GLY:HA3	1:A:253:GLY:HA2	1.89	0.54
1:G:13:LEU:O	1:G:15:LYS:NZ	2.38	0.54
1:H:87:MET:SD	3:P:277:THR:HG23	2.48	0.54
1:E:156:GLU:HG2	1:E:229:THR:HG22	1.90	0.54
1:B:10:GLN:NE2	1:B:166:SER:OG	2.40	0.54
1:B:90:ASP:OD1	2:J:279:ILE:HG13	2.08	0.54
1:E:14:PRO:HB2	1:E:17:TRP:CD1	2.43	0.54
1:F:85:LEU:HD22	3:O:260:PHE:HE2	1.73	0.54
1:C:23:GLY:HA3	1:C:253:GLY:HA2	1.90	0.54
1:E:47:VAL:HG22	1:E:141:ARG:HA	1.89	0.53
1:G:29:ARG:NH2	1:G:61:ARG:HD3	2.23	0.53
1:G:32:TRP:O	1:G:36:LEU:N	2.35	0.53
2:K:265:VAL:HG13	2:K:284:ALA:HB1	1.91	0.53
1:F:208:SER:HB3	1:F:251:LYS:HB3	1.91	0.53
1:E:204:THR:HG21	4:L:301:PEE:H14	1.91	0.53
1:A:79:ASP:OD1	1:A:80:THR:N	2.41	0.53
1:F:75:VAL:HG13	1:F:76:ARG:HG2	1.90	0.53
1:G:99:ILE:N	1:G:131:GLY:O	2.40	0.53
1:C:85:LEU:O	2:K:283:LEU:N	2.39	0.53
2:I:265:VAL:HG13	2:I:284:ALA:HB1	1.91	0.53
1:G:145:ARG:NH1	1:G:246:GLU:OE2	2.34	0.53
1:D:81:ASP:HB3	1:D:84:VAL:HG22	1.91	0.53
1:A:3:ASN:O	1:A:7:LEU:HD23	2.09	0.53
1:A:114:LEU:HD21	1:A:128:PHE:HB2	1.91	0.53
1:B:222:GLU:OE1	1:B:226:LYS:NZ	2.33	0.53
1:G:4:SER:HA	1:G:7:LEU:HB2	1.90	0.53
2:M:268:VAL:HG12	2:M:270:GLN:H	1.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:MET:HG3	1:A:241:LEU:HG	1.92	0.52
1:D:153:ILE:HG13	1:D:240:ALA:HB2	1.89	0.52
1:A:74:GLU:HA	1:C:242:LEU:HD13	1.91	0.52
1:F:134:VAL:HG13	3:O:258:ASN:HB2	1.90	0.52
1:F:148:MET:HG3	1:F:241:LEU:HG	1.90	0.52
1:A:204:THR:HG21	4:I:301:PEE:H14	1.92	0.52
1:H:47:VAL:HA	1:H:141:ARG:HG3	1.91	0.52
1:B:208:SER:HB3	1:B:251:LYS:HB3	1.92	0.52
1:E:114:LEU:HD21	1:E:128:PHE:HB2	1.92	0.51
1:E:168:ASN:ND2	1:E:170:LEU:HB2	2.24	0.51
1:H:148:MET:HE3	1:H:149:PRO:HD2	1.92	0.51
1:C:148:MET:HG3	1:C:241:LEU:HG	1.93	0.51
1:B:88:PRO:HD3	2:J:283:LEU:HD22	1.93	0.51
1:E:38:ILE:HG12	1:E:66:PHE:CD2	2.45	0.51
1:F:114:LEU:HD21	1:F:128:PHE:HB2	1.92	0.51
1:E:208:SER:HB3	1:E:251:LYS:HB3	1.93	0.51
1:H:-1:ASP:OD1	1:H:-1:ASP:N	2.38	0.51
1:H:9:LEU:O	1:H:13:LEU:HB2	2.11	0.51
1:G:112:TYR:HE2	1:G:117:LEU:HD22	1.76	0.51
1:G:151:ASN:O	1:G:191:THR:OG1	2.21	0.51
1:C:153:ILE:HG13	1:C:240:ALA:HB2	1.92	0.51
1:H:13:LEU:HD12	1:H:14:PRO:HD2	1.93	0.51
1:B:210:GLU:HB3	1:B:216:THR:HA	1.93	0.50
1:A:85:LEU:N	2:I:284:ALA:O	2.27	0.50
1:A:165:PHE:HB2	1:A:178:LEU:HD11	1.93	0.50
1:F:143:TYR:CZ	1:F:145:ARG:HB2	2.46	0.50
1:H:143:TYR:CZ	1:H:145:ARG:HB2	2.45	0.50
1:D:23:GLY:HA3	1:D:253:GLY:HA2	1.92	0.50
1:D:148:MET:HG3	1:D:241:LEU:HG	1.93	0.50
1:A:14:PRO:HB2	1:A:17:TRP:CD1	2.46	0.50
1:G:114:LEU:HD21	1:G:128:PHE:HB2	1.94	0.50
1:A:136:THR:HB	2:I:256:VAL:HG12	1.93	0.50
1:G:168:ASN:N	4:M:301:PEE:O2P	2.44	0.50
1:A:78:ILE:HG12	1:A:244:GLY:HA3	1.94	0.50
1:D:102:ASP:OD1	1:D:114:LEU:N	2.43	0.50
1:H:179:PHE:HE1	3:P:255:THR:HG21	1.76	0.50
1:D:101:GLU:HG2	1:D:130:ASN:OD1	2.12	0.50
1:F:136:THR:HB	3:O:256:VAL:HG22	1.94	0.49
1:A:41:PHE:CD2	1:A:63:PHE:HZ	2.22	0.49
1:E:137:TYR:OH	4:L:301:PEE:O2P	2.24	0.49
1:F:35:LYS:NZ	1:F:39:ASP:OD2	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:87:MET:HG3	2:L:281:GLN:HB2	1.95	0.49
1:G:29:ARG:HA	1:G:62:THR:HG22	1.94	0.49
1:G:106:GLN:HB2	1:G:112:TYR:CE1	2.47	0.49
1:A:203:ALA:HB3	1:A:206:VAL:HB	1.93	0.49
1:E:9:LEU:HA	1:E:12:ILE:HG22	1.93	0.49
1:H:161:PRO:HD2	1:H:225:ILE:HG12	1.94	0.49
2:L:261:ALA:HB3	2:L:264:LYS:HB2	1.95	0.49
1:F:15:LYS:H	1:F:15:LYS:HD3	1.78	0.49
1:F:3:ASN:HB3	1:F:170:LEU:HD11	1.93	0.49
1:F:161:PRO:HD2	1:F:225:ILE:HG12	1.94	0.49
1:D:191:THR:HG23	1:D:193:PHE:N	2.28	0.48
1:F:145:ARG:NH1	1:F:246:GLU:OE2	2.40	0.48
1:A:11:TYR:CE1	1:E:17:TRP:HH2	2.30	0.48
1:D:101:GLU:O	1:D:102:ASP:HB3	2.13	0.48
1:E:153:ILE:HG13	1:E:240:ALA:HB2	1.96	0.48
1:F:144:HIS:O	1:F:249:ARG:HB2	2.13	0.48
1:F:147:HIS:CD2	3:O:279:ILE:HD13	2.47	0.48
1:F:170:LEU:O	1:F:174:ASN:ND2	2.46	0.48
1:H:10:GLN:OE1	1:H:166:SER:N	2.36	0.48
1:H:151:ASN:OD1	1:H:243:LYS:N	2.47	0.48
1:B:177:ASN:HB3	1:B:180:ALA:HB3	1.95	0.48
3:O:268:VAL:HG12	3:O:270:GLN:H	1.78	0.48
1:D:137:TYR:CE1	1:D:167:VAL:HG13	2.47	0.48
1:E:87:MET:SD	2:L:277:THR:HG23	2.53	0.48
1:G:2:LEU:O	1:G:5:PHE:HB2	2.13	0.48
1:A:99:ILE:HG23	1:A:131:GLY:H	1.78	0.48
1:G:179:PHE:CE1	2:M:255:THR:HG21	2.49	0.48
1:H:193:PHE:HE2	3:P:286:SER:HB2	1.79	0.48
1:F:158:ILE:HB	1:F:186:ILE:HB	1.96	0.48
1:D:90:ASP:HA	3:N:277:THR:O	2.13	0.48
1:E:104:ILE:HD11	1:E:114:LEU:HD13	1.95	0.48
1:F:73:ASP:OD1	1:F:74:GLU:N	2.46	0.48
1:G:29:ARG:HH21	1:G:61:ARG:HD3	1.77	0.47
1:G:143:TYR:CZ	1:G:145:ARG:HB2	2.49	0.47
1:D:198:GLN:OE1	3:N:258:ASN:ND2	2.41	0.47
1:E:99:ILE:O	1:E:130:ASN:HA	2.13	0.47
1:A:167:VAL:HG12	4:I:301:PEE:O4P	2.14	0.47
1:H:35:LYS:NZ	1:H:39:ASP:OD2	2.46	0.47
1:G:19:THR:OG1	1:G:204:THR:O	2.26	0.47
1:A:203:ALA:HB2	4:I:301:PEE:H12	1.97	0.47
3:O:264:LYS:HD2	3:O:264:LYS:HA	1.65	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:GLY:HA2	2:J:262:PRO:CA	2.44	0.47
1:F:179:PHE:CE1	3:O:255:THR:HG21	2.49	0.47
1:H:193:PHE:CE2	3:P:286:SER:HB2	2.50	0.47
1:G:130:ASN:O	2:M:262:PRO:HB3	2.15	0.47
1:C:9:LEU:HD11	1:C:13:LEU:HD13	1.96	0.46
1:E:9:LEU:HD23	1:E:10:GLN:H	1.80	0.46
1:E:143:TYR:CZ	1:E:145:ARG:HB2	2.51	0.46
1:H:114:LEU:HD23	1:H:129:ARG:HG3	1.96	0.46
1:D:72:ARG:O	1:D:75:VAL:HG12	2.15	0.46
1:E:138:LEU:O	4:L:301:PEE:H11	2.15	0.46
1:F:210:GLU:HG2	1:F:249:ARG:HG2	1.98	0.46
1:A:153:ILE:HG13	1:A:240:ALA:HB2	1.97	0.46
1:B:99:ILE:HG13	1:B:103:LYS:O	2.15	0.46
1:C:81:ASP:HB3	1:C:84:VAL:HG22	1.97	0.46
1:E:99:ILE:HG23	1:E:104:ILE:HG12	1.98	0.46
1:E:151:ASN:OD1	1:E:243:LYS:N	2.49	0.46
1:B:3:ASN:HA	1:B:6:LYS:HD3	1.97	0.46
1:C:168:ASN:N	4:K:301:PEE:O2P	2.49	0.46
1:G:60:TYR:HD1	1:G:65:GLU:HG2	1.81	0.46
1:C:143:TYR:CZ	1:C:145:ARG:HB2	2.50	0.45
3:O:258:ASN:HB3	3:O:260:PHE:CE1	2.50	0.45
1:G:38:ILE:O	1:G:42:VAL:HG23	2.16	0.45
1:G:131:GLY:HA2	2:M:262:PRO:HA	1.97	0.45
1:H:106:GLN:HB2	1:H:112:TYR:CD1	2.51	0.45
1:B:151:ASN:OD1	1:B:243:LYS:N	2.39	0.45
1:C:86:VAL:HA	2:K:282:PRO:HA	1.99	0.45
1:F:137:TYR:CE1	1:F:167:VAL:HG13	2.51	0.45
1:A:145:ARG:HG2	1:A:249:ARG:HB3	1.99	0.45
1:A:197:ALA:HB3	2:I:259:LEU:HB2	1.98	0.45
1:C:79:ASP:OD1	1:C:80:THR:N	2.49	0.45
1:E:85:LEU:HD23	1:E:149:PRO:HB2	1.99	0.45
1:F:45:TYR:HB3	1:F:141:ARG:HB2	1.98	0.45
1:G:166:SER:OG	4:M:301:PEE:H13	2.16	0.45
1:D:193:PHE:HB3	3:N:265:VAL:HG23	1.98	0.45
1:F:4:SER:O	1:F:8:SER:N	2.38	0.45
1:H:95:GLN:H	1:H:135:THR:HB	1.80	0.45
1:F:225:ILE:HD12	1:H:119:ALA:HB1	1.98	0.45
1:H:81:ASP:HB3	1:H:84:VAL:HG22	1.97	0.45
1:D:106:GLN:HB2	1:D:112:TYR:CD1	2.51	0.45
1:B:137:TYR:OH	4:J:301:PEE:O1P	2.17	0.45
1:D:53:GLN:HB3	1:D:70:PRO:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:THR:HG23	1:D:217:ILE:HD11	2.00	0.44
1:F:14:PRO:HB2	1:F:17:TRP:HB2	1.99	0.44
1:C:78:ILE:HD12	1:C:78:ILE:H	1.82	0.44
1:D:90:ASP:OD2	1:D:143:TYR:HB3	2.18	0.44
1:F:52:ALA:HA	1:F:71:LEU:HD23	1.98	0.44
1:F:64:ASN:O	1:F:68:VAL:HG22	2.17	0.44
1:E:89:ALA:HB2	1:E:146:VAL:HA	1.97	0.44
1:E:233:GLY:O	1:E:235:ASN:ND2	2.49	0.44
1:G:75:VAL:HG13	1:G:76:ARG:HG2	1.98	0.44
1:G:77:PRO:O	2:M:280:GLY:HA3	2.18	0.44
1:G:167:VAL:HG12	1:G:167:VAL:O	2.17	0.44
1:D:193:PHE:CE1	3:N:286:SER:HB2	2.53	0.44
1:H:211:THR:HG23	1:H:217:ILE:HD11	1.98	0.44
1:B:89:ALA:HB2	1:B:146:VAL:HA	1.98	0.44
1:D:197:ALA:HB3	3:N:259:LEU:HB2	2.00	0.44
1:G:198:GLN:NE2	2:M:258:ASN:OD1	2.46	0.44
1:H:40:LEU:HD23	1:H:40:LEU:HA	1.87	0.44
1:F:106:GLN:HB2	1:F:112:TYR:CD1	2.53	0.44
1:F:195:PRO:HD2	3:O:261:ALA:HB2	1.98	0.44
1:B:81:ASP:O	1:B:243:LYS:NZ	2.51	0.44
1:E:148:MET:HG3	1:E:241:LEU:HG	2.00	0.44
1:H:143:TYR:CE2	1:H:145:ARG:HB2	2.53	0.44
1:B:106:GLN:HB2	1:B:112:TYR:CE2	2.52	0.44
1:E:52:ALA:HA	1:E:71:LEU:HA	2.00	0.44
1:A:86:VAL:HG22	2:I:279:ILE:O	2.18	0.44
1:E:165:PHE:O	1:E:204:THR:N	2.51	0.43
1:A:99:ILE:HG23	1:A:131:GLY:N	2.32	0.43
1:B:143:TYR:CZ	1:B:145:ARG:HB2	2.53	0.43
1:C:180:ALA:O	1:C:184:ARG:NH2	2.50	0.43
1:G:145:ARG:HA	1:G:249:ARG:HA	2.00	0.43
1:A:219:PRO:HG3	1:A:221:ARG:NH1	2.33	0.43
1:A:134:VAL:HG12	2:I:258:ASN:HB2	2.00	0.43
1:B:168:ASN:OD1	1:B:168:ASN:N	2.49	0.43
1:B:235:ASN:H	1:B:238:SER:HB3	1.83	0.43
3:N:273:SER:HB2	3:N:274:LEU:HD22	1.99	0.43
1:H:206:VAL:CG1	1:H:250:PHE:HB3	2.49	0.43
1:D:64:ASN:OD1	1:D:252:LEU:HB2	2.17	0.43
1:A:26:ALA:HB1	1:A:63:PHE:HB3	2.01	0.43
1:D:38:ILE:HG12	1:D:66:PHE:CD2	2.53	0.43
1:G:186:ILE:HG23	1:G:199:ILE:HG12	2.00	0.43
3:N:267:LEU:HA	3:N:284:ALA:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:106:GLN:HB2	1:E:112:TYR:CD2	2.54	0.43
1:G:193:PHE:CE2	2:M:286:SER:HB2	2.54	0.43
1:A:81:ASP:O	1:A:84:VAL:HG22	2.19	0.43
1:E:72:ARG:O	1:E:75:VAL:HG12	2.18	0.43
1:A:42:VAL:HA	1:A:47:VAL:HG23	1.99	0.42
1:H:199:ILE:HD12	3:P:257:ILE:HD11	2.01	0.42
1:A:143:TYR:CZ	1:A:145:ARG:HB2	2.55	0.42
1:F:13:LEU:O	1:F:15:LYS:HD3	2.18	0.42
1:F:86:VAL:HB	1:F:243:LYS:HE3	2.01	0.42
1:E:33:LEU:H	1:E:33:LEU:HD12	1.84	0.42
1:B:166:SER:N	1:B:204:THR:HB	2.35	0.42
1:B:179:PHE:CE1	2:J:255:THR:HG21	2.54	0.42
1:E:165:PHE:H	1:E:182:ASN:ND2	2.11	0.42
2:I:273:SER:HB2	2:I:274:LEU:HD22	2.02	0.42
1:B:51:GLU:HG2	1:B:75:VAL:HG21	2.01	0.42
1:H:168:ASN:ND2	1:H:170:LEU:HB2	2.35	0.42
1:C:87:MET:HA	2:K:283:LEU:HD22	2.01	0.42
1:G:88:PRO:HG3	2:M:258:ASN:HD21	1.85	0.42
1:H:93:ILE:HD12	3:P:275:SER:HB2	2.02	0.42
1:H:146:VAL:HG21	3:P:256:VAL:HG11	2.02	0.42
1:A:159:TYR:OH	1:A:183:GLU:OE2	2.31	0.42
1:A:162:GLY:HA3	1:A:181:ARG:O	2.20	0.42
1:C:9:LEU:O	1:C:13:LEU:HB2	2.20	0.42
1:H:34:THR:O	1:H:38:ILE:HG13	2.20	0.42
1:A:205:ILE:HG21	1:A:205:ILE:HD13	1.75	0.42
1:B:96:LEU:HD12	2:J:267:LEU:HD22	2.02	0.42
1:B:148:MET:HG3	1:B:241:LEU:HG	2.00	0.42
1:D:10:GLN:HE21	1:D:10:GLN:HB3	1.65	0.42
1:D:159:TYR:HB3	1:D:226:LYS:HB2	2.01	0.42
1:D:168:ASN:OD1	1:D:168:ASN:N	2.53	0.42
1:G:219:PRO:HG3	1:G:221:ARG:HH12	1.85	0.42
1:H:211:THR:HA	1:H:248:GLY:HA2	2.01	0.42
1:B:1:MET:O	1:B:4:SER:HB2	2.20	0.41
1:B:87:MET:HA	2:J:283:LEU:HD22	2.00	0.41
1:H:13:LEU:O	1:H:15:LYS:NZ	2.48	0.41
1:H:177:ASN:HB3	1:H:180:ALA:HB3	2.01	0.41
1:A:15:LYS:O	1:A:19:THR:OG1	2.37	0.41
1:C:42:VAL:HG13	1:C:47:VAL:HB	2.02	0.41
1:F:4:SER:HA	1:F:7:LEU:HB2	2.01	0.41
1:H:137:TYR:CE1	1:H:167:VAL:HG13	2.55	0.41
1:A:76:ARG:HA	1:A:76:ARG:HD3	1.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:VAL:HG13	1:C:76:ARG:HG2	2.02	0.41
1:E:19:THR:HG22	1:E:205:ILE:HA	2.01	0.41
1:F:1:MET:O	1:F:4:SER:HB2	2.21	0.41
1:F:205:ILE:HD13	1:F:205:ILE:HA	1.91	0.41
1:A:7:LEU:HD23	1:A:7:LEU:H	1.85	0.41
1:B:195:PRO:HD2	2:J:261:ALA:HB2	2.02	0.41
1:C:199:ILE:O	2:K:256:VAL:HA	2.20	0.41
2:K:268:VAL:HB	2:K:271:LEU:HD12	2.02	0.41
1:D:241:LEU:HD11	1:D:247:MET:HB3	2.02	0.41
1:F:38:ILE:HG12	1:F:66:PHE:CD2	2.55	0.41
3:O:258:ASN:HB3	3:O:260:PHE:HE1	1.86	0.41
1:F:84:VAL:HA	3:O:285:VAL:HA	2.02	0.41
1:B:17:TRP:HA	1:B:17:TRP:CE3	2.55	0.41
1:C:29:ARG:HB3	1:C:62:THR:HG22	2.02	0.41
1:H:104:ILE:HD11	1:H:114:LEU:HD13	2.03	0.41
1:H:148:MET:HG3	1:H:241:LEU:HG	2.02	0.41
1:H:205:ILE:HD13	1:H:205:ILE:HA	1.91	0.41
1:D:12:ILE:HD12	1:D:12:ILE:HA	1.93	0.41
1:E:167:VAL:HA	1:E:178:LEU:HD21	2.03	0.41
1:G:99:ILE:HD11	1:G:128:PHE:O	2.21	0.41
1:G:250:PHE:CD2	2:M:254:PPI:H33	2.54	0.41
1:C:150:CYS:HA	1:C:243:LYS:HB2	2.03	0.41
1:D:3:ASN:O	1:D:7:LEU:HD13	2.21	0.41
1:D:86:VAL:HG12	1:D:149:PRO:O	2.21	0.41
1:E:143:TYR:OH	1:E:145:ARG:HB2	2.21	0.41
4:L:301:PEE:O1P	4:L:301:PEE:H12	2.20	0.41
1:F:233:GLY:HA2	1:F:238:SER:OG	2.21	0.41
1:G:4:SER:O	1:G:8:SER:OG	2.31	0.41
1:G:155:ARG:CZ	1:G:232:ALA:HA	2.51	0.41
1:H:96:LEU:HD21	3:P:272:GLU:HA	2.03	0.41
1:G:91:GLY:HA2	1:G:142:ASP:OD2	2.21	0.40
1:H:210:GLU:HB3	1:H:216:THR:HA	2.03	0.40
1:A:99:ILE:O	1:A:130:ASN:HA	2.21	0.40
1:C:137:TYR:HE1	1:C:167:VAL:HG13	1.86	0.40
1:G:204:THR:HG21	4:M:301:PEE:H14	2.03	0.40
1:B:147:HIS:CD2	2:J:279:ILE:HD13	2.56	0.40
1:G:122:TYR:HA	1:G:125:ALA:HB3	2.03	0.40
1:G:159:TYR:OH	1:G:183:GLU:OE2	2.31	0.40
1:A:-1:ASP:HA	1:A:2:LEU:HB3	2.04	0.40
1:D:33:LEU:O	1:D:37:VAL:HG23	2.21	0.40
1:E:166:SER:O	4:L:301:PEE:O1P	2.40	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	253/267 (95%)	228 (90%)	25 (10%)	0	100	100
1	B	253/267 (95%)	229 (90%)	24 (10%)	0	100	100
1	C	253/267 (95%)	230 (91%)	23 (9%)	0	100	100
1	D	253/267 (95%)	228 (90%)	25 (10%)	0	100	100
1	E	253/267 (95%)	233 (92%)	19 (8%)	1 (0%)	34	71
1	F	253/267 (95%)	229 (90%)	23 (9%)	1 (0%)	34	71
1	G	253/267 (95%)	232 (92%)	20 (8%)	1 (0%)	34	71
1	H	253/267 (95%)	230 (91%)	23 (9%)	0	100	100
2	I	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
2	J	32/36 (89%)	28 (88%)	4 (12%)	0	100	100
2	K	32/36 (89%)	29 (91%)	3 (9%)	0	100	100
2	L	32/36 (89%)	28 (88%)	4 (12%)	0	100	100
2	M	31/36 (86%)	29 (94%)	2 (6%)	0	100	100
3	N	32/36 (89%)	29 (91%)	3 (9%)	0	100	100
3	O	32/36 (89%)	29 (91%)	3 (9%)	0	100	100
3	P	32/36 (89%)	29 (91%)	3 (9%)	0	100	100
All	All	2279/2424 (94%)	2071 (91%)	205 (9%)	3 (0%)	51	83

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	167	VAL
1	F	88	PRO
1	G	206	VAL

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	216/229 (94%)	210 (97%)	6 (3%)	43	72
1	B	215/229 (94%)	207 (96%)	8 (4%)	34	66
1	C	209/229 (91%)	203 (97%)	6 (3%)	42	72
1	D	214/229 (93%)	205 (96%)	9 (4%)	30	63
1	E	211/229 (92%)	201 (95%)	10 (5%)	26	61
1	F	213/229 (93%)	207 (97%)	6 (3%)	43	72
1	G	215/229 (94%)	209 (97%)	6 (3%)	43	72
1	H	209/229 (91%)	205 (98%)	4 (2%)	57	80
2	I	27/31 (87%)	26 (96%)	1 (4%)	34	66
2	J	27/31 (87%)	27 (100%)	0	100	100
2	K	27/31 (87%)	27 (100%)	0	100	100
2	L	27/31 (87%)	27 (100%)	0	100	100
2	M	27/31 (87%)	27 (100%)	0	100	100
3	N	27/31 (87%)	27 (100%)	0	100	100
3	O	27/31 (87%)	27 (100%)	0	100	100
3	P	26/31 (84%)	26 (100%)	0	100	100
All	All	1917/2080 (92%)	1861 (97%)	56 (3%)	42	72

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	20	ARG
1	A	29	ARG
1	A	41	PHE
1	A	113	SER
1	A	236	ASP
2	I	264	LYS
1	B	20	ARG

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Mol	Chain	Res	Type
1	B	33	LEU
1	B	69	ARG
1	B	73	ASP
1	B	164	LEU
1	B	218	THR
1	B	236	ASP
1	B	238	SER
1	C	20	ARG
1	C	53	GLN
1	C	69	ARG
1	C	113	SER
1	C	139	SER
1	C	182	ASN
1	D	1	MET
1	D	13	LEU
1	D	20	ARG
1	D	32	TRP
1	D	66	PHE
1	D	69	ARG
1	D	143	TYR
1	D	173	GLN
1	D	236	ASP
1	E	-1	ASP
1	E	5	PHE
1	E	11	TYR
1	E	13	LEU
1	E	20	ARG
1	E	66	PHE
1	E	69	ARG
1	E	122	TYR
1	E	143	TYR
1	E	236	ASP
1	F	5	PHE
1	F	15	LYS
1	F	20	ARG
1	F	66	PHE
1	F	143	TYR
1	F	182	ASN
1	G	20	ARG
1	G	29	ARG
1	G	69	ARG
1	G	73	ASP

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Mol	Chain	Res	Type
1	G	151	ASN
1	G	236	ASP
1	H	1	MET
1	H	20	ARG
1	H	69	ARG
1	H	96	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	147	HIS
1	E	168	ASN
1	G	198	GLN
2	M	258	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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
4	PEE	M	301	2	22,22,50	1.39	2 (9%)	25,27,55	1.11	2 (8%)
4	PEE	I	301	2	22,22,50	1.41	2 (9%)	25,27,55	1.38	4 (16%)
4	PEE	L	301	2	22,22,50	1.45	2 (9%)	25,27,55	1.33	3 (12%)
4	PEE	J	301	2	22,22,50	1.41	2 (9%)	25,27,55	1.23	3 (12%)
4	PEE	K	301	2	22,22,50	1.41	2 (9%)	25,27,55	1.14	3 (12%)

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
4	PEE	M	301	2	-	11/26/26/54	-
4	PEE	I	301	2	-	13/26/26/54	-
4	PEE	L	301	2	-	11/26/26/54	-
4	PEE	J	301	2	-	11/26/26/54	-
4	PEE	K	301	2	-	11/26/26/54	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	301	PEE	O2-C10	4.67	1.47	1.34
4	I	301	PEE	O2-C10	4.45	1.46	1.34
4	J	301	PEE	O2-C10	4.42	1.46	1.34
4	K	301	PEE	O2-C10	4.42	1.46	1.34
4	M	301	PEE	O2-C10	4.34	1.46	1.34
4	L	301	PEE	O3-C30	4.33	1.46	1.33
4	K	301	PEE	O3-C30	4.23	1.45	1.33
4	J	301	PEE	O3-C30	4.23	1.45	1.33
4	M	301	PEE	O3-C30	4.20	1.45	1.33
4	I	301	PEE	O3-C30	4.00	1.45	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	301	PEE	O2-C10-C11	4.56	121.33	111.50
4	J	301	PEE	O2-C10-C11	3.83	119.76	111.50
4	L	301	PEE	O2-C10-C11	3.82	119.74	111.50
4	M	301	PEE	O2-C10-C11	3.34	118.70	111.50
4	K	301	PEE	O2-C10-C11	3.25	118.51	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	301	PEE	C3-C2-C1	-3.01	104.66	111.79
4	L	301	PEE	O3-C30-C31	2.80	120.70	111.91
4	J	301	PEE	O3-C30-C31	2.53	119.85	111.91
4	K	301	PEE	O3-C30-C31	2.51	119.80	111.91
4	I	301	PEE	C3-O3-C30	2.40	126.01	117.12
4	M	301	PEE	O3-C30-C31	2.28	119.06	111.91
4	I	301	PEE	O3-C30-C31	2.14	118.63	111.91
4	J	301	PEE	O3-C30-O5	-2.11	118.26	123.59
4	K	301	PEE	O3-C30-O5	-2.06	118.40	123.59
4	I	301	PEE	O3-C3-C2	2.01	114.28	108.43

There are no chirality outliers.

All (57) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	I	301	PEE	O2-C2-C3-O3
4	I	301	PEE	C1-O3P-P-O2P
4	I	301	PEE	C1-O3P-P-O1P
4	I	301	PEE	C4-O4P-P-O2P
4	J	301	PEE	C11-C10-O2-C2
4	M	301	PEE	C2-C1-O3P-P
4	L	301	PEE	C2-C3-O3-C30
4	J	301	PEE	O5-C30-O3-C3
4	K	301	PEE	O5-C30-O3-C3
4	L	301	PEE	O5-C30-O3-C3
4	J	301	PEE	O4-C10-O2-C2
4	J	301	PEE	C31-C30-O3-C3
4	L	301	PEE	C31-C30-O3-C3
4	K	301	PEE	C31-C30-O3-C3
4	K	301	PEE	C2-C3-O3-C30
4	J	301	PEE	C2-C1-O3P-P
4	I	301	PEE	C11-C10-O2-C2
4	I	301	PEE	C1-O3P-P-O4P
4	I	301	PEE	C4-O4P-P-O3P
4	M	301	PEE	C1-O3P-P-O4P
4	I	301	PEE	O4-C10-O2-C2
4	M	301	PEE	O4-C10-O2-C2
4	M	301	PEE	C11-C10-O2-C2
4	K	301	PEE	C11-C10-O2-C2
4	K	301	PEE	O4-C10-O2-C2
4	L	301	PEE	C4-O4P-P-O3P
4	I	301	PEE	C2-C1-O3P-P

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Mol	Chain	Res	Type	Atoms
4	M	301	PEE	C1-C2-O2-C10
4	I	301	PEE	C1-C2-C3-O3
4	J	301	PEE	C1-C2-O2-C10
4	M	301	PEE	O2-C2-C3-O3
4	I	301	PEE	C30-C31-C32-C33
4	K	301	PEE	C10-C11-C12-C13
4	L	301	PEE	C4-O4P-P-O2P
4	M	301	PEE	C1-O3P-P-O1P
4	I	301	PEE	C2-C3-O3-C30
4	J	301	PEE	O4P-C4-C5-N
4	L	301	PEE	C5-C4-O4P-P
4	M	301	PEE	C1-C2-C3-O3
4	K	301	PEE	C3-C2-O2-C10
4	J	301	PEE	C1-O3P-P-O4P
4	K	301	PEE	C2-C1-O3P-P
4	J	301	PEE	C10-C11-C12-C13
4	K	301	PEE	O2-C2-C3-O3
4	L	301	PEE	C1-C2-O2-C10
4	K	301	PEE	O4P-C4-C5-N
4	M	301	PEE	O4P-C4-C5-N
4	L	301	PEE	C10-C11-C12-C13
4	L	301	PEE	C3-C2-O2-C10
4	J	301	PEE	C1-O3P-P-O1P
4	J	301	PEE	C4-O4P-P-O1P
4	K	301	PEE	C4-O4P-P-O1P
4	L	301	PEE	C1-O3P-P-O1P
4	L	301	PEE	C4-O4P-P-O1P
4	I	301	PEE	C5-C4-O4P-P
4	M	301	PEE	C5-C4-O4P-P
4	M	301	PEE	O3-C30-C31-C32

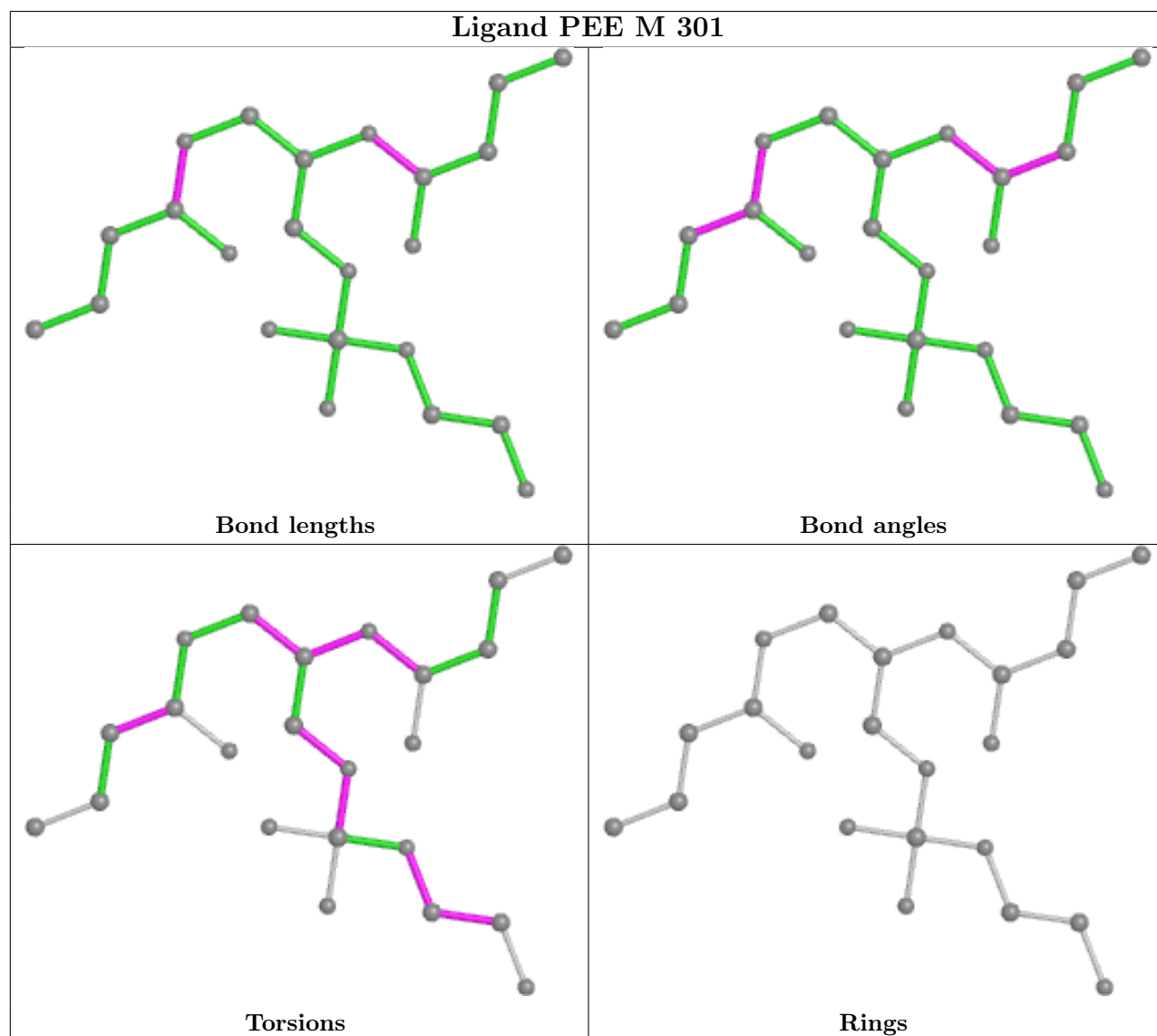
There are no ring outliers.

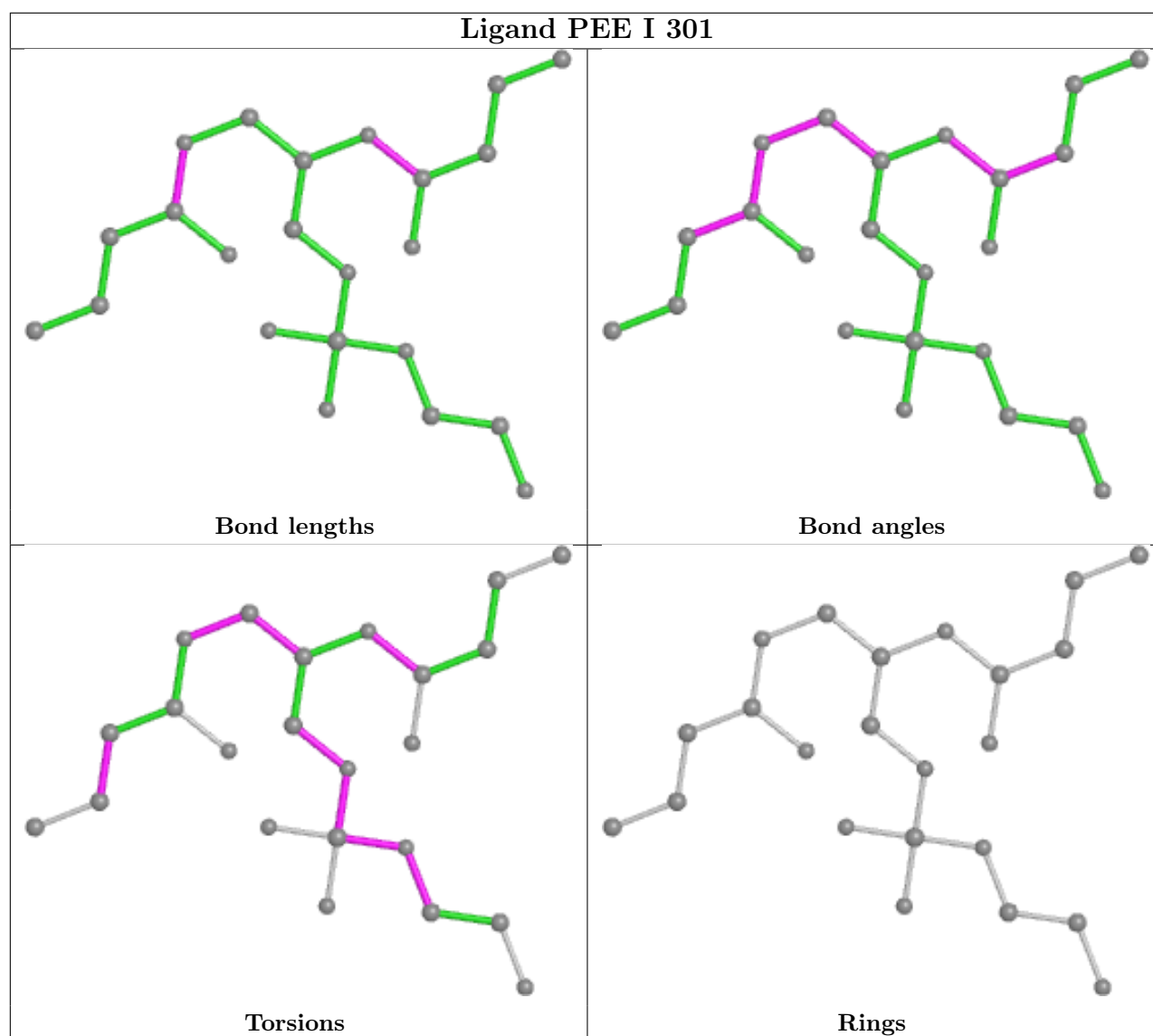
5 monomers are involved in 19 short contacts:

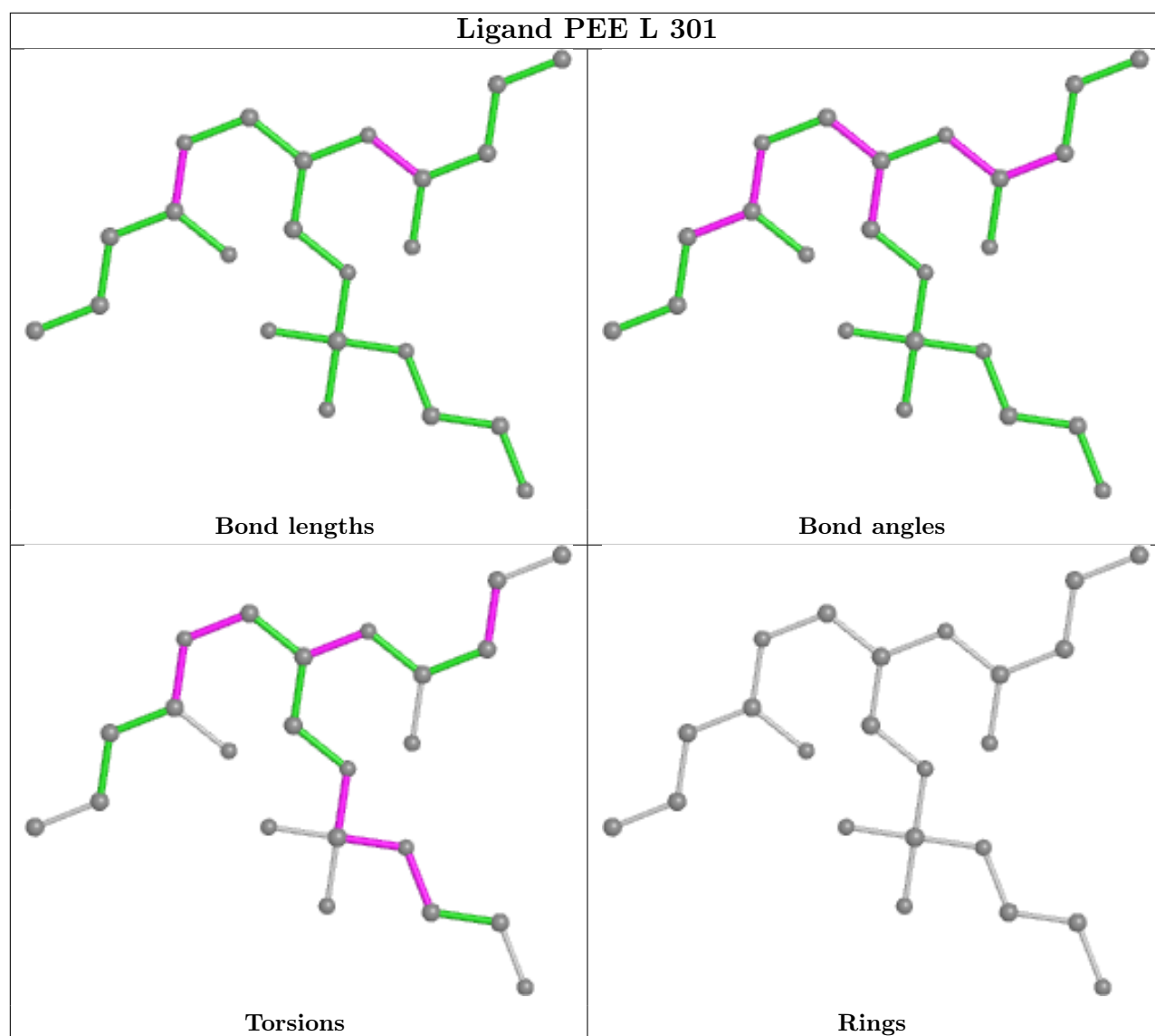
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	M	301	PEE	3	0
4	I	301	PEE	7	0
4	L	301	PEE	5	0
4	J	301	PEE	1	0
4	K	301	PEE	3	0

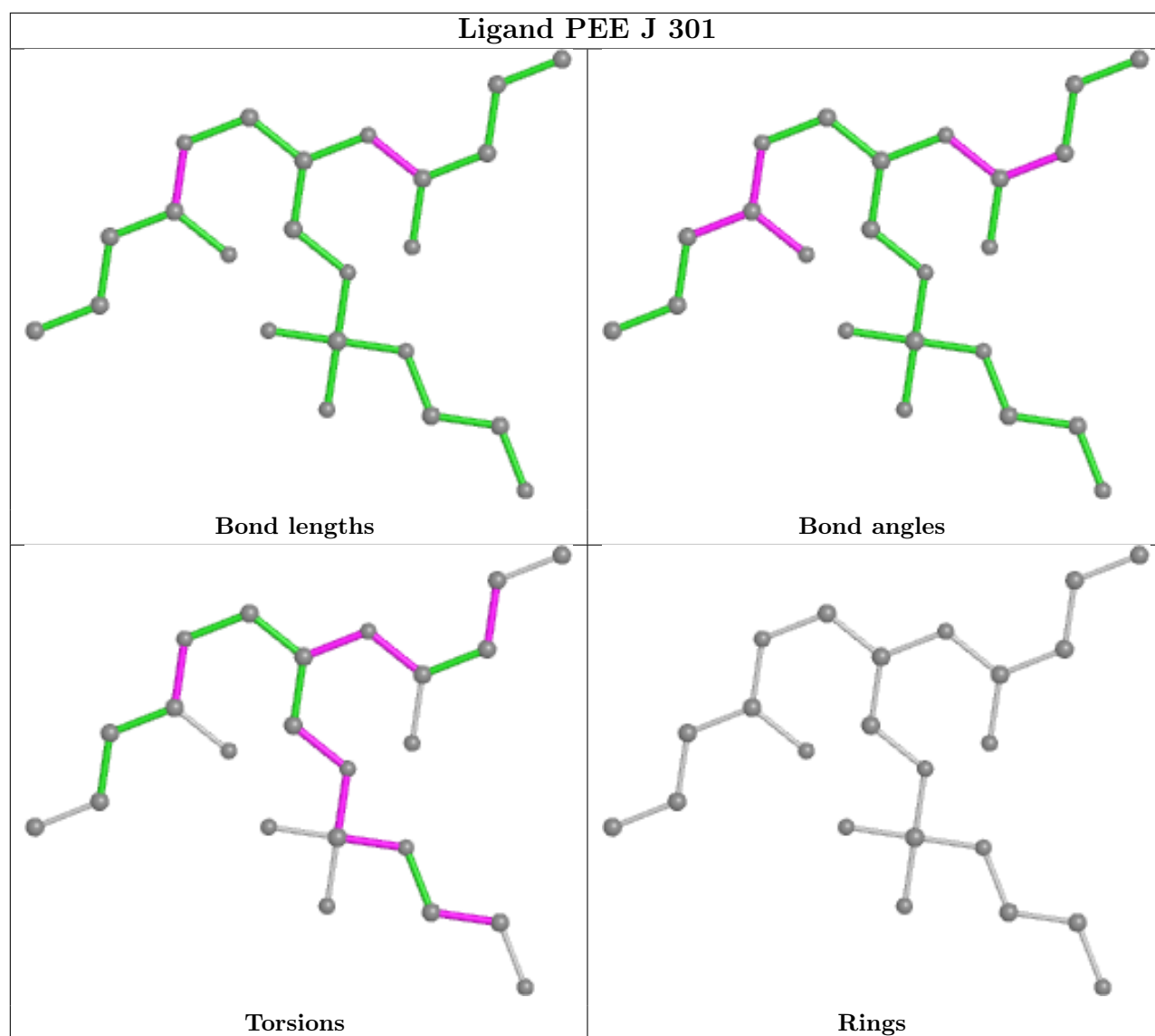
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

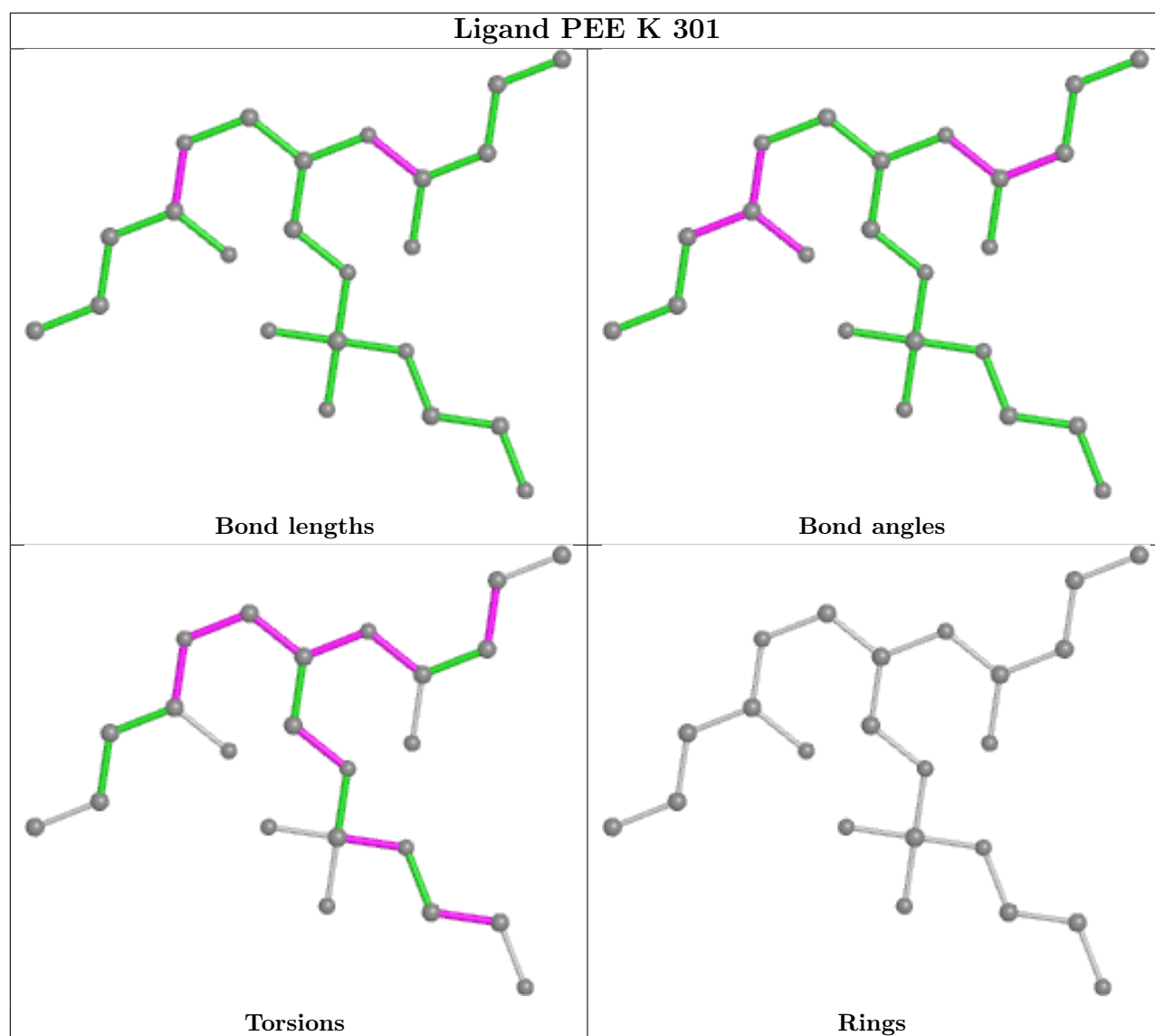
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











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	255/267 (95%)	-0.33	0 100 100	106, 129, 165, 190	0
1	B	255/267 (95%)	-0.35	0 100 100	106, 137, 170, 196	0
1	C	255/267 (95%)	-0.29	0 100 100	115, 140, 177, 198	0
1	D	255/267 (95%)	-0.31	1 (0%) 92 86	108, 144, 181, 209	0
1	E	255/267 (95%)	-0.30	4 (1%) 72 57	119, 156, 184, 194	0
1	F	255/267 (95%)	-0.34	0 100 100	131, 160, 184, 195	0
1	G	255/267 (95%)	-0.19	3 (1%) 79 66	121, 162, 187, 210	0
1	H	255/267 (95%)	0.20	17 (6%) 17 10	152, 189, 214, 226	0
2	I	34/36 (94%)	-0.13	0 100 100	104, 127, 146, 157	0
2	J	34/36 (94%)	-0.15	0 100 100	118, 145, 160, 173	0
2	K	34/36 (94%)	-0.47	0 100 100	128, 157, 182, 192	0
2	L	34/36 (94%)	-0.53	0 100 100	136, 168, 184, 195	0
2	M	33/36 (91%)	-0.01	0 100 100	150, 179, 193, 196	0
3	N	34/36 (94%)	-0.17	1 (2%) 51 35	125, 155, 170, 172	0
3	O	34/36 (94%)	0.04	1 (2%) 51 35	143, 166, 176, 183	0
3	P	34/36 (94%)	0.40	5 (14%) 2 1	180, 205, 218, 222	0
All	All	2311/2424 (95%)	-0.23	32 (1%) 75 61	104, 153, 199, 226	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	196	MET	6.6
1	H	187	CYS	4.8
1	H	189	PHE	4.8
3	P	260	PHE	4.6
1	H	195	PRO	4.3

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Mol	Chain	Res	Type	RSRZ
1	H	149	PRO	3.6
1	H	188	LEU	3.4
1	H	198	GLN	3.3
1	H	148	MET	3.3
3	P	259	LEU	3.2
3	P	288	GLU	3.1
1	H	197	ALA	3.1
1	H	61	ARG	3.1
3	P	258	ASN	2.8
1	H	86	VAL	2.8
1	G	179	PHE	2.7
1	H	85	LEU	2.6
3	O	259	LEU	2.6
1	G	209	ILE	2.6
1	D	135	THR	2.3
3	N	282	PRO	2.3
1	E	187	CYS	2.3
1	H	63	PHE	2.2
1	E	157	MET	2.2
1	H	194	GLY	2.1
1	E	221	ARG	2.1
1	E	202	GLY	2.1
1	H	154	LEU	2.1
1	G	252	LEU	2.1
1	H	66	PHE	2.1
3	P	279	ILE	2.0
1	H	57	THR	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 monosaccharides 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,

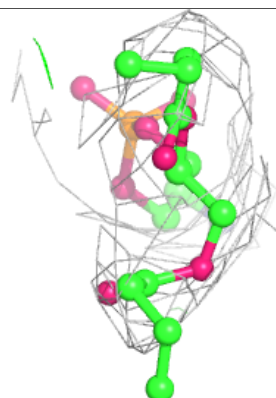
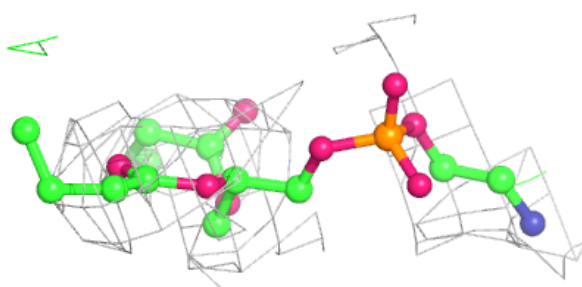
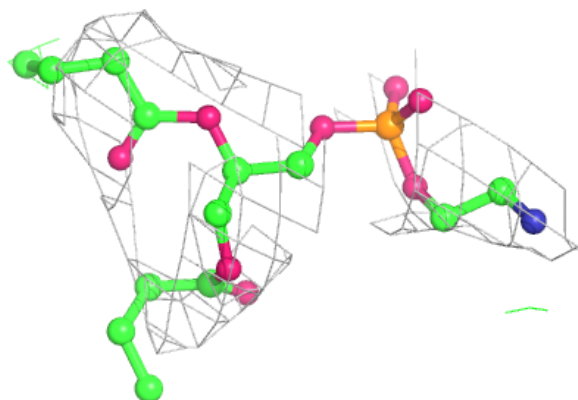
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PEE	L	301	23/51	0.89	0.27	109,142,164,167	0
4	PEE	I	301	23/51	0.90	0.29	106,142,148,155	0
4	PEE	M	301	23/51	0.90	0.26	129,148,161,168	0
4	PEE	J	301	23/51	0.93	0.25	110,138,150,158	0
4	PEE	K	301	23/51	0.93	0.31	105,130,143,156	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

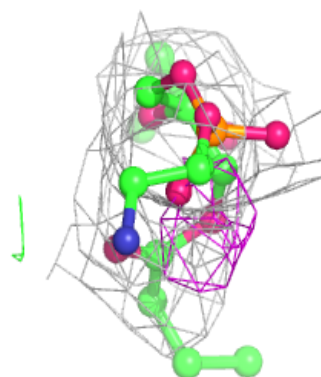
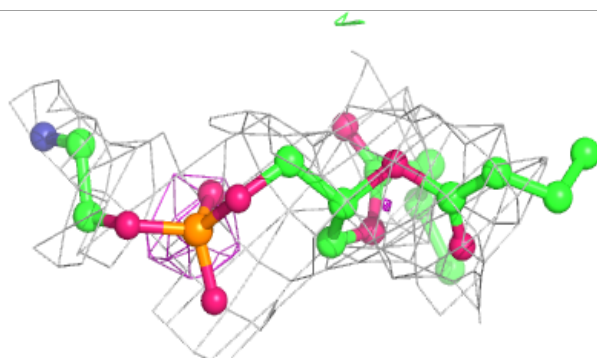
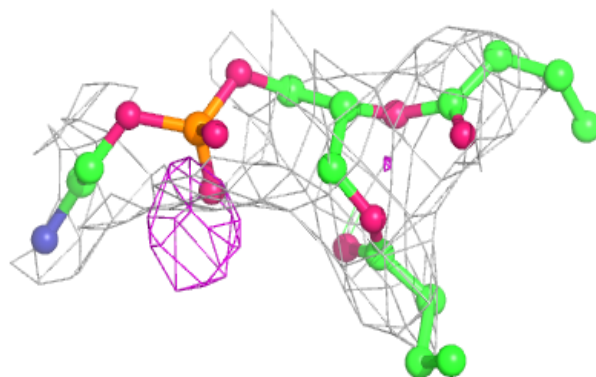
Electron density around PEE L 301:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

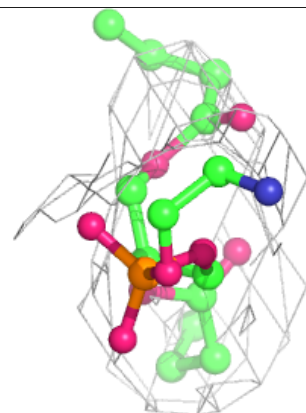
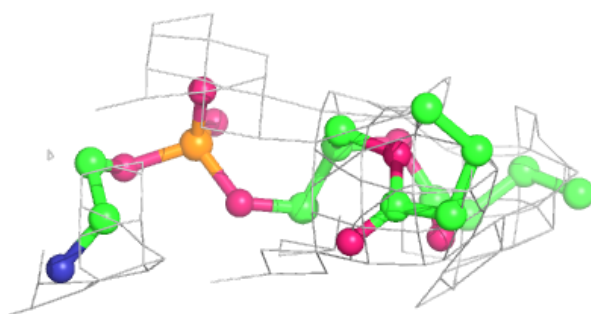
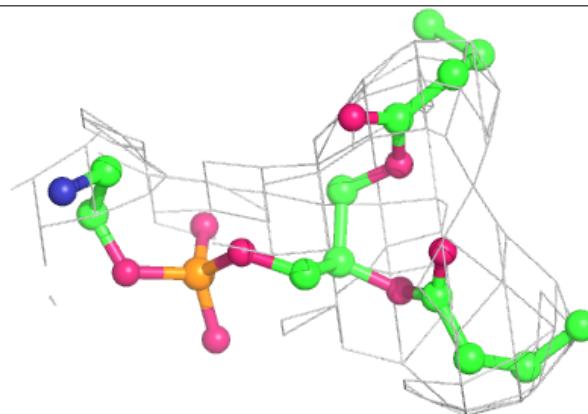


Electron density around PEE I 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

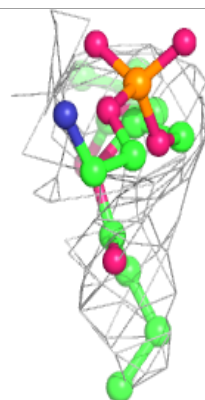
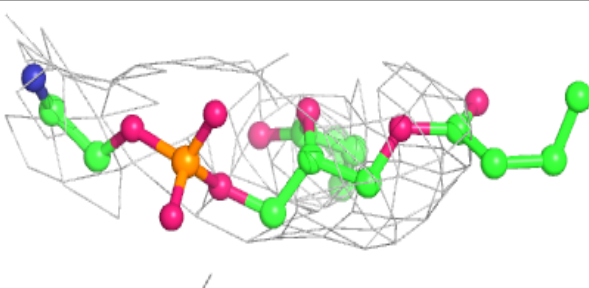
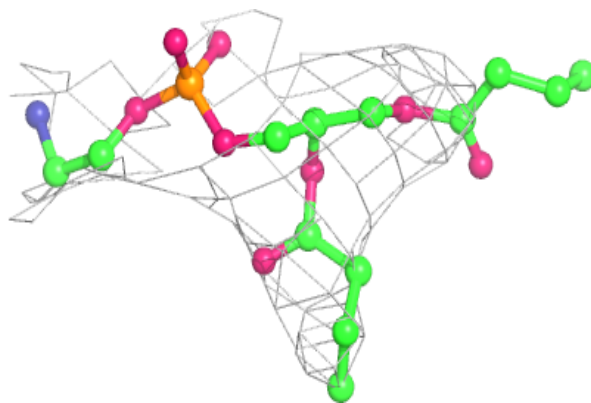
**Electron density around PEE M 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

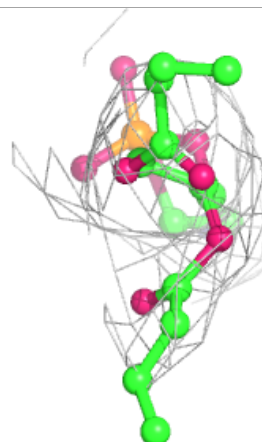
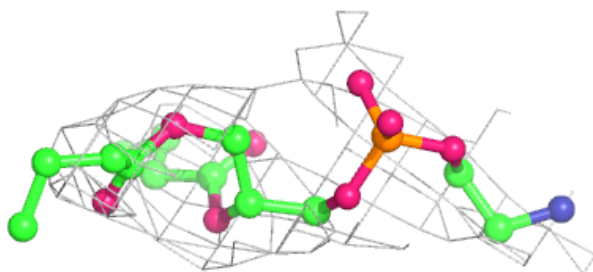
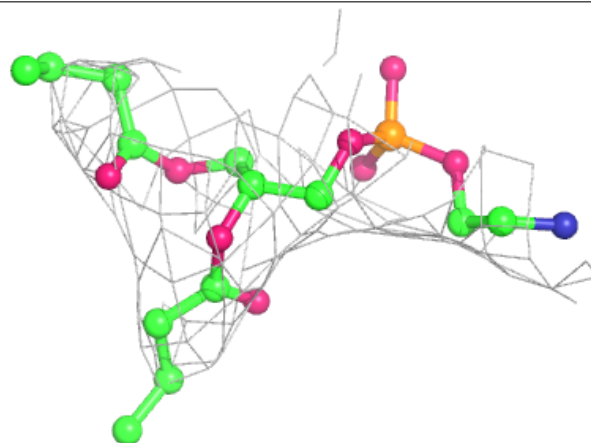


Electron density around PEE J 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around PEE K 301:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.