



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 12:12 am GMT

PDB ID : 5E9S
Title : Crystal structure of substrate-bound glutamate transporter homologue GltTk
Authors : Guskov, A.; Slotboom, D.J.
Deposited on : 2015-10-15
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

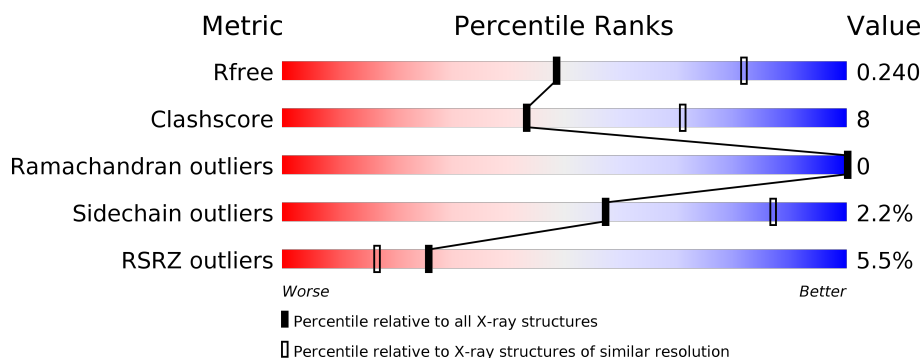
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	438	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>21%</div> <div>• •</div> </div> </div>
1	B	438	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>17%</div> <div>•</div> </div> </div>
1	C	438	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>•</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	DMU	A	505	-	-	-	X
4	DMU	B	505	-	-	-	X
4	DMU	C	505	-	-	-	X
5	1PE	A	506	-	-	-	X
5	1PE	B	506	-	-	-	X
5	1PE	B	507	-	-	-	X
5	1PE	C	506	-	-	-	X
6	PEG	A	508	-	-	-	X
6	PEG	A	510	-	-	-	X
6	PEG	A	512	-	-	-	X
6	PEG	A	516	-	-	-	X
6	PEG	B	509	-	-	-	X
6	PEG	B	510	-	-	-	X
6	PEG	B	511	-	-	-	X
6	PEG	B	516	-	-	-	X
6	PEG	B	518	-	-	-	X
6	PEG	B	519	-	-	-	X
6	PEG	C	510	-	-	-	X
6	PEG	C	513	-	-	-	X
6	PEG	C	514	-	-	-	X
6	PEG	C	516	-	-	-	X
7	PGE	A	520	-	-	-	X
7	PGE	A	521	-	-	-	X
7	PGE	C	518	-	-	-	X
8	P6G	A	522	-	-	-	X
8	P6G	B	525	-	-	-	X
8	P6G	C	519	-	-	-	X
8	P6G	C	520	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10213 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 Proton/glutamate symporter, SDF family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	1	0
			3196	2108	518	553	17			
1	B	427	Total	C	N	O	S	0	0	0
			3188	2103	517	552	16			
1	C	426	Total	C	N	O	S	0	1	0
			3186	2102	515	552	17			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	431	HIS	-	expression tag	UNP Q5JID0
A	432	HIS	-	expression tag	UNP Q5JID0
A	433	HIS	-	expression tag	UNP Q5JID0
A	434	HIS	-	expression tag	UNP Q5JID0
A	435	HIS	-	expression tag	UNP Q5JID0
A	436	HIS	-	expression tag	UNP Q5JID0
A	437	HIS	-	expression tag	UNP Q5JID0
A	438	HIS	-	expression tag	UNP Q5JID0
B	431	HIS	-	expression tag	UNP Q5JID0
B	432	HIS	-	expression tag	UNP Q5JID0
B	433	HIS	-	expression tag	UNP Q5JID0
B	434	HIS	-	expression tag	UNP Q5JID0
B	435	HIS	-	expression tag	UNP Q5JID0
B	436	HIS	-	expression tag	UNP Q5JID0
B	437	HIS	-	expression tag	UNP Q5JID0
B	438	HIS	-	expression tag	UNP Q5JID0
C	431	HIS	-	expression tag	UNP Q5JID0
C	432	HIS	-	expression tag	UNP Q5JID0
C	433	HIS	-	expression tag	UNP Q5JID0
C	434	HIS	-	expression tag	UNP Q5JID0
C	435	HIS	-	expression tag	UNP Q5JID0
C	436	HIS	-	expression tag	UNP Q5JID0
C	437	HIS	-	expression tag	UNP Q5JID0

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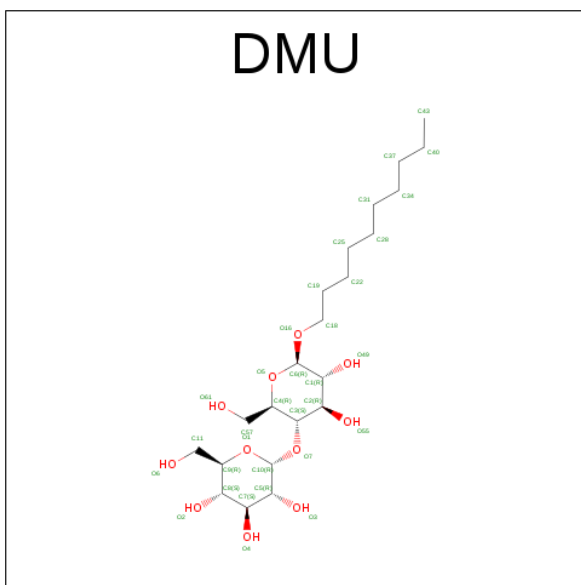
Chain	Residue	Modelled	Actual	Comment	Reference
C	438	HIS	-	expression tag	UNP Q5JID0

- # ASP
-
- The diagram shows the chemical structure of Aspartic acid (ASP). The central carbon atom is bonded to a hydrogen atom (H), an amino group (NH₂), a carboxyl group (COOH), and a side chain. The side chain consists of a methylene group (CH₂) and another carboxyl group (COOH). The atoms are color-coded: carbon (grey), oxygen (red), hydrogen (white), nitrogen (blue), and sulfur (yellow). The side chain is labeled with 'CA(S)' for the central carbon, 'CB' for the methylene carbon, and 'CG' for the carboxyl carbon. The carboxyl groups are labeled with 'OD1' and 'OD2' for the hydroxyl oxygens, and 'OXT' for the side chain hydroxyl oxygen. The amino group is labeled with 'N'.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 9	C 4	N 1	O 4	0	0
2	B	1	Total 9	C 4	N 1	O 4	0	0
2	C	1	Total 9	C 4	N 1	O 4	0	0

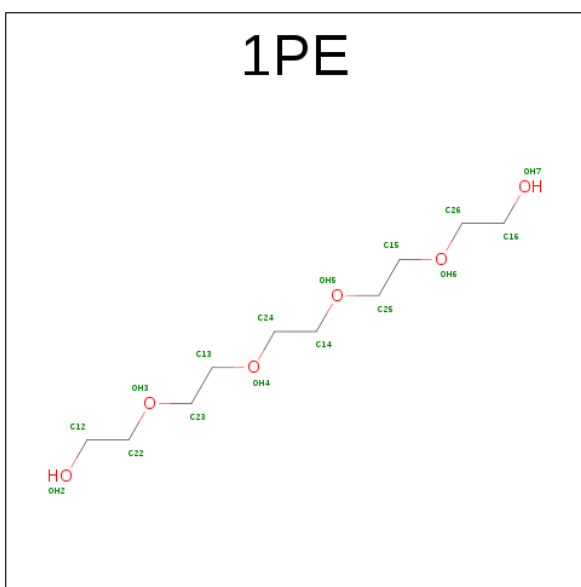
- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 3 | B | 3 | Total Na
3 3 | 0 | 0 |
| 3 | A | 3 | Total Na
3 3 | 0 | 0 |
| 3 | C | 3 | Total Na
3 3 | 0 | 0 |

- 
- WORLD WIDE
PDB
PROTEIN DATA BANK



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 33	C 22	O 11	0	0
4	B	1	Total 33	C 22	O 11	0	0
4	C	1	Total 33	C 22	O 11	0	0

- Molecule 5 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $\text{C}_{10}\text{H}_{22}\text{O}_6$).



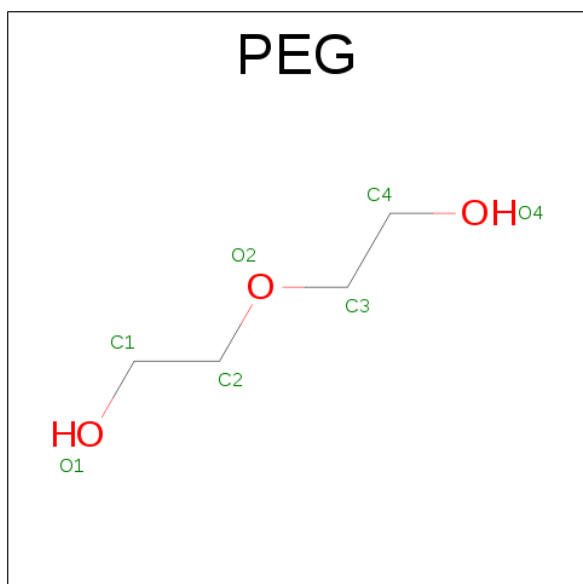
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			16	10	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	C	1	Total	C	O	0	0
			16	10	6		
5	C	1	Total	C	O	0	0
			16	10	6		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		

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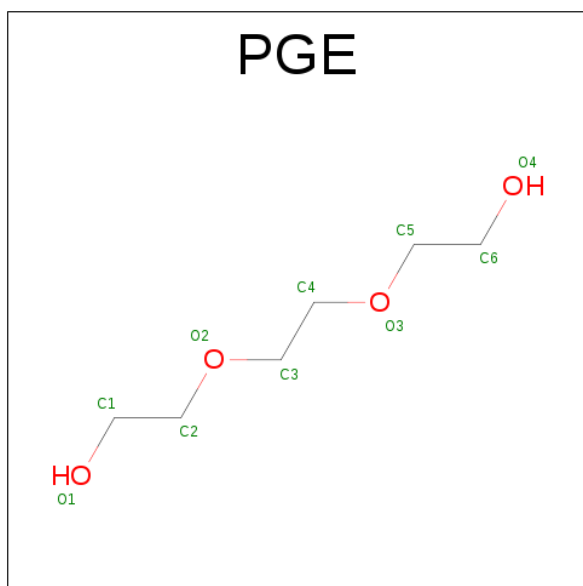
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



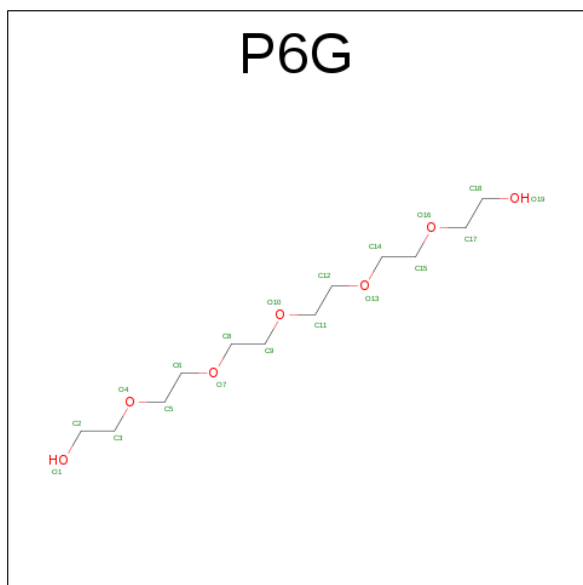
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			10	6	4		
7	A	1	Total	C	O	0	0
			10	6	4		
7	A	1	Total	C	O	0	0
			10	6	4		
7	A	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			10	6	4		
7	C	1	Total	C	O	0	0
			10	6	4		
7	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 8 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).

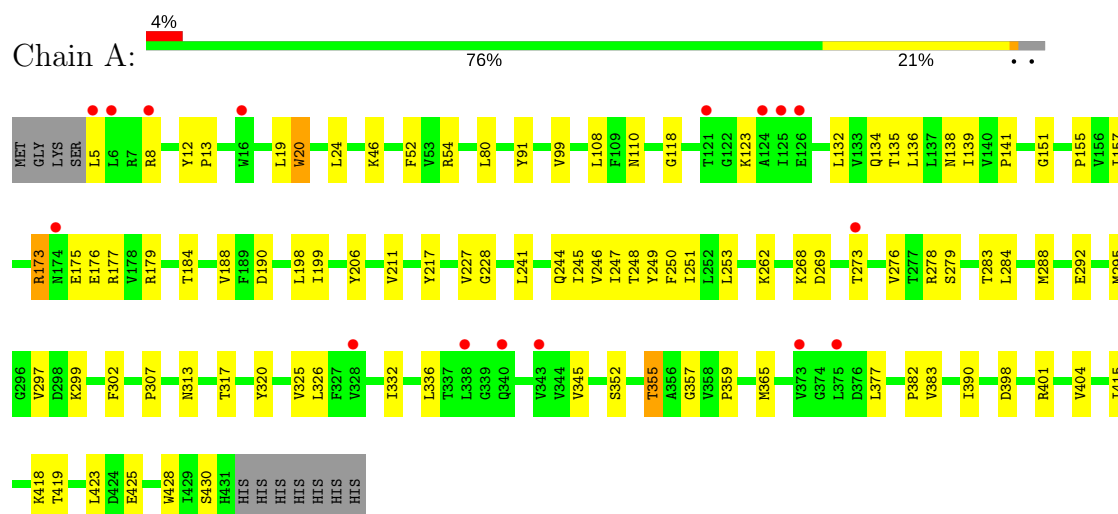


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			19	12	7		
8	B	1	Total	C	O	0	0
			19	12	7		
8	B	1	Total	C	O	0	0
			19	12	7		
8	C	1	Total	C	O	0	0
			19	12	7		
8	C	1	Total	C	O	0	0
			19	12	7		

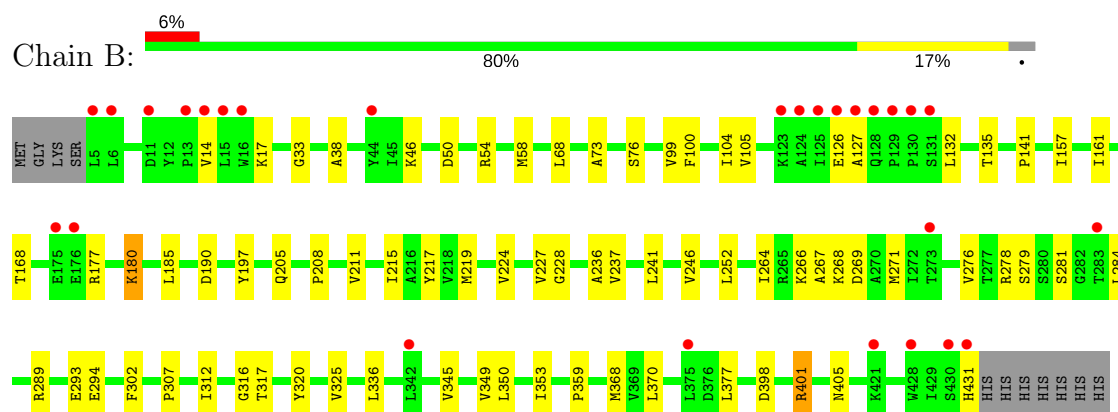
3 Residue-property plots [i](#)

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

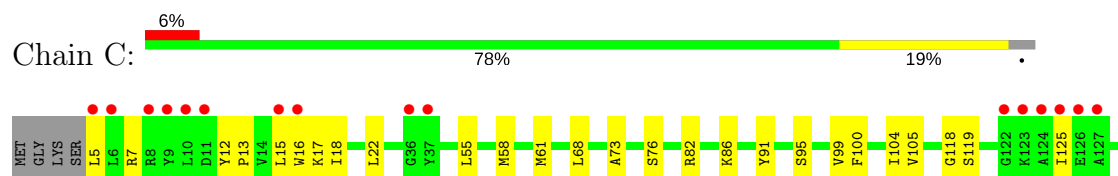
- Molecule 1: Proton/glutamate symporter, SDF family

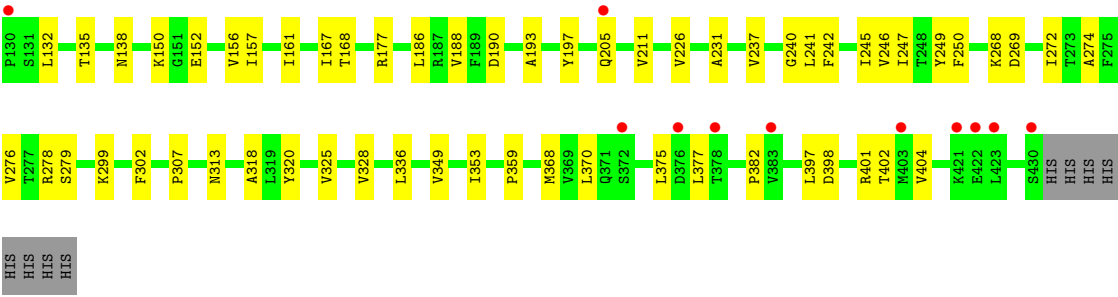


- Molecule 1: Proton/glutamate symporter, SDF family



- Molecule 1: Proton/glutamate symporter, SDF family





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.52Å 117.52Å 310.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.36 – 2.80 48.36 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.36-2.80) 97.4 (48.36-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.213 , 0.242 0.210 , 0.240	Depositor DCC
R_{free} test set	2010 reflections (3.32%)	DCC
Wilson B-factor (Å ²)	83.7	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 65.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10213	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PGE, NA, 1PE, DMU, P6G, PEG

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	0.43	0/3256	0.64	0/4431
1	B	0.42	0/3248	0.61	0/4421
1	C	0.40	0/3245	0.63	0/4416
All	All	0.41	0/9749	0.63	0/13268

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	118	GLY	Peptide

5.2 Too-close contacts [i](#)

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	3196	0	3402	64	0
1	B	3188	0	3394	45	0
1	C	3186	0	3395	53	0
2	A	9	0	3	2	0
2	B	9	0	3	3	0
2	C	9	0	3	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
4	A	33	0	42	0	0
4	B	33	0	42	1	0
4	C	33	0	42	2	0
5	A	32	0	44	2	0
5	B	48	0	66	4	0
5	C	32	0	44	0	0
6	A	70	0	100	1	0
6	B	98	0	140	3	0
6	C	63	0	90	5	0
7	A	40	0	56	2	0
7	B	10	0	14	0	0
7	C	20	0	28	0	0
8	A	19	0	26	0	0
8	B	38	0	52	2	0
8	C	38	0	52	1	0
All	All	10213	0	11038	163	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 (163) 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)
5:B:508:1PE:H241	6:B:522:PEG:H22	1.63	0.81
1:A:357:GLY:H	7:A:520:PGE:H12	1.47	0.79
1:A:110:ASN:HB2	5:A:507:1PE:H251	1.65	0.78
1:A:177:ARG:NH1	1:B:190:ASP:OD2	2.20	0.74
1:C:241:LEU:HB3	1:C:404:VAL:HG21	1.72	0.70
1:B:157:ILE:HD11	1:B:307:PRO:HB2	1.73	0.69
1:C:12:TYR:HD1	1:C:17:LYS:HD3	1.57	0.69
1:C:157:ILE:HD11	1:C:307:PRO:HB2	1.74	0.69
1:A:246:VAL:HG12	1:A:247:ILE:HD12	1.74	0.69
1:A:157:ILE:HD11	1:A:307:PRO:HB2	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:LEU:HB3	1:C:16:TRP:HD1	1.58	0.68
1:B:14:VAL:HG13	1:B:17:LYS:HE2	1.77	0.66
1:C:119:SER:HA	1:C:382:PRO:HG2	1.78	0.65
1:A:190:ASP:OD2	1:C:177:ARG:NH1	2.30	0.65
1:B:99:VAL:HG11	1:B:345:VAL:HA	1.79	0.64
1:A:355:THR:HG22	1:A:365:MET:HG3	1.79	0.64
1:A:262:LYS:HE2	1:A:428:TRP:O	2.00	0.62
1:A:5:LEU:HD23	1:A:8:ARG:HD3	1.80	0.61
1:C:246:VAL:HG23	1:C:247:ILE:HD12	1.82	0.61
5:B:507:1PE:H221	6:B:509:PEG:H32	1.82	0.60
1:C:274:ALA:HB1	1:C:402:THR:HG22	1.83	0.59
1:C:325:VAL:HG12	1:C:336:LEU:HD11	1.82	0.59
1:C:167:ILE:HG21	1:C:186:LEU:HD13	1.86	0.58
1:C:55:LEU:O	1:C:58:MET:HG2	2.03	0.58
1:A:279:SER:HB2	1:A:359:PRO:HD3	1.85	0.57
1:B:17:LYS:HE3	1:B:208:PRO:HG2	1.85	0.57
1:C:246:VAL:O	1:C:250:PHE:HB2	2.05	0.57
1:C:237:VAL:O	1:C:241:LEU:HG	2.05	0.56
1:B:177:ARG:NH1	1:C:190:ASP:OD2	2.38	0.56
1:A:418:LYS:NZ	1:A:425:GLU:OE2	2.40	0.55
1:B:325:VAL:HG22	1:B:370:LEU:HD12	1.89	0.55
1:A:284:LEU:HD22	1:A:288:MET:HE2	1.88	0.55
1:A:325:VAL:HG12	1:A:336:LEU:HD11	1.87	0.55
1:C:12:TYR:CD1	1:C:17:LYS:HD3	2.39	0.55
1:B:205:GLN:NE2	4:B:505:DMU:O49	2.36	0.55
1:B:278:ARG:HD2	1:B:398:ASP:HB3	1.89	0.55
1:A:20:TRP:HA	1:A:20:TRP:CE3	2.41	0.54
1:A:269:ASP:O	1:A:273:THR:HG23	2.08	0.54
1:A:262:LYS:NZ	1:A:430:SER:OG	2.38	0.54
1:C:197:TYR:HH	1:C:302:PHE:HE1	1.53	0.54
1:C:325:VAL:HG22	1:C:370:LEU:HD12	1.90	0.54
1:B:73:ALA:O	1:B:168:THR:HG21	2.09	0.53
1:A:20:TRP:HA	1:A:20:TRP:HE3	1.73	0.53
1:B:266:LYS:HB3	1:B:294:GLU:HB3	1.91	0.52
1:C:278:ARG:HD2	1:C:398:ASP:HB3	1.90	0.52
1:A:99:VAL:HG11	1:A:345:VAL:HA	1.92	0.51
1:A:401:ARG:NH1	2:A:501:ASP:OD1	2.43	0.51
1:C:99:VAL:HG22	1:C:318:ALA:HB1	1.93	0.51
1:A:12:TYR:CG	1:A:13:PRO:HD2	2.46	0.51
1:A:52:PHE:HB2	1:A:206:TYR:CE2	2.45	0.51
1:B:126:GLU:HB2	1:B:127:ALA:HA	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:ARG:HD2	1:A:398:ASP:HB3	1.93	0.50
1:C:226:VAL:HG13	1:C:231:ALA:HA	1.93	0.50
2:B:501:ASP:OD1	2:B:501:ASP:N	2.42	0.50
1:B:267:ALA:O	1:B:271:MET:HG3	2.12	0.50
1:A:184:THR:O	1:A:188:VAL:HG23	2.12	0.49
1:A:295:MET:HB2	1:A:297:VAL:HG23	1.94	0.49
1:A:54:ARG:NH2	1:C:138:ASN:HA	2.28	0.49
1:A:211:VAL:HG22	1:A:276:VAL:HG21	1.94	0.49
1:B:325:VAL:HG12	1:B:336:LEU:HD11	1.94	0.48
1:A:141:PRO:HD2	1:B:58:MET:HE3	1.94	0.48
1:A:132:LEU:O	1:A:135:THR:HB	2.14	0.48
1:A:245:ILE:HA	1:A:249:TYR:CD2	2.49	0.48
1:B:317:THR:OG1	2:B:501:ASP:OXT	2.30	0.48
1:A:108:LEU:HA	5:A:507:1PE:H242	1.95	0.48
1:B:279:SER:HB2	1:B:359:PRO:HD3	1.95	0.47
1:A:199:ILE:HG23	6:C:510:PEG:H11	1.96	0.47
1:A:423:LEU:HD11	1:A:428:TRP:HE1	1.79	0.47
1:A:246:VAL:O	1:A:250:PHE:HB2	2.15	0.47
6:C:508:PEG:H22	6:C:509:PEG:H21	1.95	0.47
1:C:73:ALA:O	1:C:168:THR:OG1	2.32	0.47
1:A:139:ILE:O	1:A:155:PRO:HA	2.15	0.47
1:A:317:THR:OG1	2:A:501:ASP:OD2	2.29	0.47
1:C:105:VAL:HG21	1:C:240:GLY:HA2	1.97	0.47
1:C:132:LEU:O	1:C:135:THR:N	2.47	0.47
1:C:242:PHE:O	1:C:246:VAL:HG22	2.14	0.47
1:B:197:TYR:HH	1:B:302:PHE:HE1	1.63	0.47
1:C:22:LEU:HD22	1:C:272:ILE:HG12	1.95	0.47
1:C:150:LYS:HD3	1:C:152:GLU:OE2	2.14	0.47
1:C:91:TYR:CG	1:C:313:ASN:HB2	2.50	0.46
1:A:151:GLY:HA3	7:A:521:PGE:H4	1.98	0.46
1:B:281:SER:O	1:B:284:LEU:HB2	2.15	0.46
1:A:118:GLY:O	1:A:382:PRO:HG2	2.15	0.46
1:A:227:VAL:HG13	1:A:228:GLY:N	2.31	0.45
1:B:132:LEU:O	1:B:135:THR:HB	2.16	0.45
1:B:353:ILE:HG21	5:B:506:1PE:H252	1.97	0.45
1:C:68:LEU:HD21	1:C:161:ILE:HG13	1.98	0.45
1:A:173:ARG:NH2	1:A:175:GLU:OE1	2.50	0.45
1:C:5:LEU:HG	1:C:7:ARG:H	1.81	0.45
1:A:176:GLU:HG3	1:A:179:ARG:NH2	2.32	0.45
1:A:268:LYS:HG3	1:A:269:ASP:N	2.32	0.44
1:B:405:ASN:ND2	2:B:501:ASP:HB2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:LYS:HD2	1:A:217:TYR:CG	2.53	0.44
1:C:349:VAL:O	1:C:353:ILE:HG13	2.17	0.44
8:C:520:P6G:H92	8:C:520:P6G:H141	2.00	0.44
1:A:377:LEU:HD23	1:A:377:LEU:HA	1.53	0.44
1:B:211:VAL:O	1:B:215:ILE:HG22	2.18	0.44
5:B:508:1PE:H131	6:B:522:PEG:H32	1.98	0.44
1:A:123:LYS:HE3	1:A:123:LYS:HB3	1.71	0.44
1:A:273:THR:OG1	1:A:283:THR:HG23	2.16	0.44
1:A:332:ILE:HD13	1:A:383:VAL:HG22	1.99	0.44
1:C:193:ALA:O	1:C:197:TYR:HD2	2.00	0.44
1:C:245:ILE:HA	1:C:249:TYR:CD2	2.53	0.44
1:A:198:LEU:HD22	6:C:510:PEG:H41	2.00	0.44
1:C:211:VAL:HG22	1:C:276:VAL:HG21	1.99	0.44
1:B:237:VAL:O	1:B:241:LEU:HG	2.18	0.43
1:C:186:LEU:HD23	6:C:512:PEG:H32	2.00	0.43
1:C:272:ILE:O	1:C:276:VAL:HG22	2.18	0.43
1:C:268:LYS:HG3	1:C:269:ASP:N	2.33	0.43
1:C:61:MET:CE	1:C:156:VAL:HG21	2.48	0.43
1:C:279:SER:HB2	1:C:359:PRO:HD3	2.00	0.43
1:C:99:VAL:CG2	1:C:318:ALA:HB1	2.49	0.43
1:B:276:VAL:O	1:B:278:ARG:NH1	2.52	0.43
1:B:33:GLY:HA2	1:B:38:ALA:HB2	1.99	0.43
1:B:312:ILE:HG21	8:B:524:P6G:H182	2.01	0.43
1:A:317:THR:HG21	1:A:355:THR:HG21	2.01	0.43
1:A:141:PRO:O	1:B:58:MET:HB3	2.18	0.43
1:A:80:LEU:HD12	1:A:80:LEU:HA	1.72	0.42
1:C:328:VAL:HG12	1:C:375:LEU:HD13	2.01	0.42
1:C:397:LEU:O	1:C:401:ARG:HG2	2.19	0.42
1:A:247:ILE:HG23	1:A:251:ILE:HD12	2.01	0.42
8:B:525:P6G:H21	8:B:525:P6G:H52	1.84	0.42
1:A:134:GLN:HG3	1:A:138:ASN:HD21	1.84	0.42
1:A:91:TYR:CG	1:A:313:ASN:HB2	2.55	0.42
1:B:289:ARG:NE	1:B:293:GLU:OE1	2.52	0.42
1:B:316:GLY:C	1:B:401:ARG:HG3	2.40	0.42
1:C:302:PHE:CD1	1:C:302:PHE:C	2.94	0.42
6:C:515:PEG:H32	6:C:515:PEG:H11	1.78	0.42
1:C:12:TYR:CG	1:C:13:PRO:HD2	2.55	0.41
4:C:505:DMU:H35	4:C:505:DMU:H30	2.01	0.41
1:A:352:SER:HA	1:A:355:THR:HG23	2.01	0.41
1:A:326:LEU:HD23	1:A:336:LEU:HD12	2.02	0.41
1:B:105:VAL:HG11	1:B:236:ALA:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:PHE:CE2	1:B:104:ILE:HD11	2.55	0.41
1:B:180:LYS:O	1:B:180:LYS:HD2	2.20	0.41
1:A:292:GLU:HB2	1:A:302:PHE:CZ	2.55	0.41
1:B:141:PRO:HG2	1:C:58:MET:HE3	2.01	0.41
1:A:244:GLN:O	1:A:248:THR:HB	2.20	0.41
1:B:350:LEU:HD23	1:B:350:LEU:HA	1.83	0.41
1:B:268:LYS:HG3	1:B:269:ASP:N	2.35	0.41
1:C:377:LEU:HA	1:C:377:LEU:HD23	1.67	0.41
6:A:511:PEG:H31	6:A:511:PEG:H12	1.88	0.41
1:B:349:VAL:O	1:B:353:ILE:HG13	2.20	0.41
1:C:95:SER:O	1:C:99:VAL:HG23	2.20	0.41
1:A:136:LEU:O	1:A:139:ILE:HB	2.21	0.41
1:B:252:LEU:HD23	1:B:252:LEU:HA	1.96	0.41
1:B:50:ASP:O	1:B:54:ARG:HG3	2.20	0.41
1:A:241:LEU:HB3	1:A:404:VAL:HG21	2.03	0.41
1:A:415:ILE:O	1:A:419:THR:HG23	2.21	0.41
1:B:227:VAL:HG13	1:B:228:GLY:N	2.36	0.41
1:B:46:LYS:HD2	1:B:217:TYR:CD2	2.55	0.41
1:A:20:TRP:O	1:A:24:LEU:HG	2.21	0.40
1:B:215:ILE:HD11	1:B:219:MET:HE2	2.03	0.40
1:C:205:GLN:HB3	4:C:505:DMU:H13	2.03	0.40
1:A:253:LEU:HA	1:A:253:LEU:HD23	1.92	0.40
1:B:68:LEU:HD21	1:B:161:ILE:HG13	2.02	0.40
1:B:141:PRO:O	1:C:58:MET:HB3	2.21	0.40
1:B:185:LEU:HD13	1:C:188:VAL:HG13	2.03	0.40
1:C:299:LYS:HG3	1:C:302:PHE:CZ	2.56	0.40
1:A:299:LYS:HB2	1:A:299:LYS:HE2	1.89	0.40
1:A:390:ILE:HD13	1:A:390:ILE:HA	1.83	0.40
1:C:100:PHE:CE2	1:C:104:ILE:HD11	2.56	0.40
1:C:82:ARG:HG2	1:C:86:LYS:HE3	2.03	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	426/438 (97%)	418 (98%)	8 (2%)	0	100	100
1	B	425/438 (97%)	415 (98%)	10 (2%)	0	100	100
1	C	425/438 (97%)	413 (97%)	12 (3%)	0	100	100
All	All	1276/1314 (97%)	1246 (98%)	30 (2%)	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	336/345 (97%)	331 (98%)	5 (2%)	70	92
1	B	335/345 (97%)	325 (97%)	10 (3%)	46	80
1	C	335/345 (97%)	329 (98%)	6 (2%)	64	90
All	All	1006/1035 (97%)	985 (98%)	21 (2%)	57	88

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	20	TRP
1	A	173	ARG
1	A	320	TYR
1	A	355	THR
1	B	76	SER
1	B	180	LYS
1	B	224	VAL
1	B	246	VAL
1	B	264	ILE
1	B	320	TYR
1	B	368	MET
1	B	377	LEU

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Mol	Chain	Res	Type
1	B	401	ARG
1	B	431	HIS
1	C	18	ILE
1	C	76	SER
1	C	125	ILE
1	C	320	TYR
1	C	368[A]	MET
1	C	368[B]	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 67 ligands modelled in this entry, 9 are monoatomic - leaving 58 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ASP	A	501	-	1,8,8	0.41	0	1,10,10	0.55	0
4	DMU	A	505	-	34,34,34	1.70	9 (26%)	45,45,45	1.37	7 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	1PE	A	506	-	15,15,15	0.54	0	14,14,14	0.36	0
5	1PE	A	507	-	15,15,15	0.55	0	14,14,14	0.41	0
6	PEG	A	508	-	6,6,6	0.49	0	5,5,5	0.32	0
6	PEG	A	509	-	6,6,6	0.51	0	5,5,5	0.29	0
6	PEG	A	510	-	6,6,6	0.50	0	5,5,5	0.32	0
6	PEG	A	511	-	6,6,6	0.53	0	5,5,5	0.47	0
6	PEG	A	512	-	6,6,6	0.48	0	5,5,5	0.38	0
6	PEG	A	513	-	6,6,6	0.52	0	5,5,5	0.36	0
6	PEG	A	514	-	6,6,6	0.52	0	5,5,5	0.26	0
6	PEG	A	515	-	6,6,6	0.52	0	5,5,5	0.26	0
6	PEG	A	516	-	6,6,6	0.54	0	5,5,5	0.30	0
6	PEG	A	517	-	6,6,6	0.57	0	5,5,5	0.77	0
7	PGE	A	518	-	9,9,9	0.36	0	8,8,8	0.34	0
7	PGE	A	519	-	9,9,9	0.38	0	8,8,8	0.32	0
7	PGE	A	520	-	9,9,9	0.46	0	8,8,8	0.34	0
7	PGE	A	521	-	9,9,9	0.37	0	8,8,8	0.41	0
8	P6G	A	522	-	18,18,18	0.60	0	17,17,17	0.47	0
2	ASP	B	501	-	1,8,8	0.50	0	1,10,10	1.02	0
4	DMU	B	505	-	34,34,34	1.62	8 (23%)	45,45,45	1.08	5 (11%)
5	1PE	B	506	-	15,15,15	0.56	0	14,14,14	0.48	0
5	1PE	B	507	-	15,15,15	0.55	0	14,14,14	0.33	0
5	1PE	B	508	-	15,15,15	0.55	0	14,14,14	0.24	0
6	PEG	B	509	-	6,6,6	0.54	0	5,5,5	0.29	0
6	PEG	B	510	-	6,6,6	0.53	0	5,5,5	0.27	0
6	PEG	B	511	-	6,6,6	0.49	0	5,5,5	0.31	0
6	PEG	B	512	-	6,6,6	0.55	0	5,5,5	0.31	0
6	PEG	B	513	-	6,6,6	0.52	0	5,5,5	0.37	0
6	PEG	B	514	-	6,6,6	0.50	0	5,5,5	0.31	0
6	PEG	B	515	-	6,6,6	0.50	0	5,5,5	0.41	0
6	PEG	B	516	-	6,6,6	0.55	0	5,5,5	0.27	0
6	PEG	B	517	-	6,6,6	0.51	0	5,5,5	0.25	0
6	PEG	B	518	-	6,6,6	0.50	0	5,5,5	0.39	0
6	PEG	B	519	-	6,6,6	0.51	0	5,5,5	0.29	0
6	PEG	B	520	-	6,6,6	0.55	0	5,5,5	0.40	0
6	PEG	B	521	-	6,6,6	0.50	0	5,5,5	0.30	0
6	PEG	B	522	-	6,6,6	0.52	0	5,5,5	0.27	0
7	PGE	B	523	-	9,9,9	0.43	0	8,8,8	0.27	0
8	P6G	B	524	-	18,18,18	0.55	0	17,17,17	0.46	0
8	P6G	B	525	-	18,18,18	0.56	0	17,17,17	0.49	0
2	ASP	C	501	-	1,8,8	0.29	0	1,10,10	0.50	0
4	DMU	C	505	-	34,34,34	1.61	9 (26%)	45,45,45	1.11	3 (6%)
5	1PE	C	506	-	15,15,15	0.52	0	14,14,14	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	1PE	C	507	-	15,15,15	0.54	0	14,14,14	0.21	0
6	PEG	C	508	-	6,6,6	0.52	0	5,5,5	0.43	0
6	PEG	C	509	-	6,6,6	0.55	0	5,5,5	0.34	0
6	PEG	C	510	-	6,6,6	0.49	0	5,5,5	0.39	0
6	PEG	C	511	-	6,6,6	0.53	0	5,5,5	0.37	0
6	PEG	C	512	-	6,6,6	0.52	0	5,5,5	0.35	0
6	PEG	C	513	-	6,6,6	0.52	0	5,5,5	0.34	0
6	PEG	C	514	-	6,6,6	0.53	0	5,5,5	0.37	0
6	PEG	C	515	-	6,6,6	0.51	0	5,5,5	0.27	0
6	PEG	C	516	-	6,6,6	0.46	0	5,5,5	0.31	0
7	PGE	C	517	-	9,9,9	0.35	0	8,8,8	0.26	0
7	PGE	C	518	-	9,9,9	0.31	0	8,8,8	0.58	0
8	P6G	C	519	-	18,18,18	0.60	0	17,17,17	0.54	0
8	P6G	C	520	-	18,18,18	0.55	0	17,17,17	0.49	0

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	ASP	A	501	-	-	0/2/8/8	0/0/0/0
4	DMU	A	505	-	-	0/19/59/59	0/2/2/2
5	1PE	A	506	-	-	0/13/13/13	0/0/0/0
5	1PE	A	507	-	-	0/13/13/13	0/0/0/0
6	PEG	A	508	-	-	0/4/4/4	0/0/0/0
6	PEG	A	509	-	-	0/4/4/4	0/0/0/0
6	PEG	A	510	-	-	0/4/4/4	0/0/0/0
6	PEG	A	511	-	-	0/4/4/4	0/0/0/0
6	PEG	A	512	-	-	0/4/4/4	0/0/0/0
6	PEG	A	513	-	-	0/4/4/4	0/0/0/0
6	PEG	A	514	-	-	0/4/4/4	0/0/0/0
6	PEG	A	515	-	-	0/4/4/4	0/0/0/0
6	PEG	A	516	-	-	0/4/4/4	0/0/0/0
6	PEG	A	517	-	-	0/4/4/4	0/0/0/0
7	PGE	A	518	-	-	0/7/7/7	0/0/0/0
7	PGE	A	519	-	-	0/7/7/7	0/0/0/0
7	PGE	A	520	-	-	0/7/7/7	0/0/0/0
7	PGE	A	521	-	-	0/7/7/7	0/0/0/0
8	P6G	A	522	-	-	0/16/16/16	0/0/0/0
2	ASP	B	501	-	-	0/2/8/8	0/0/0/0
4	DMU	B	505	-	-	0/19/59/59	0/2/2/2
5	1PE	B	506	-	-	0/13/13/13	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	1PE	B	507	-	-	0/13/13/13	0/0/0/0
5	1PE	B	508	-	-	0/13/13/13	0/0/0/0
6	PEG	B	509	-	-	0/4/4/4	0/0/0/0
6	PEG	B	510	-	-	0/4/4/4	0/0/0/0
6	PEG	B	511	-	-	0/4/4/4	0/0/0/0
6	PEG	B	512	-	-	0/4/4/4	0/0/0/0
6	PEG	B	513	-	-	0/4/4/4	0/0/0/0
6	PEG	B	514	-	-	0/4/4/4	0/0/0/0
6	PEG	B	515	-	-	0/4/4/4	0/0/0/0
6	PEG	B	516	-	-	0/4/4/4	0/0/0/0
6	PEG	B	517	-	-	0/4/4/4	0/0/0/0
6	PEG	B	518	-	-	0/4/4/4	0/0/0/0
6	PEG	B	519	-	-	0/4/4/4	0/0/0/0
6	PEG	B	520	-	-	0/4/4/4	0/0/0/0
6	PEG	B	521	-	-	0/4/4/4	0/0/0/0
6	PEG	B	522	-	-	0/4/4/4	0/0/0/0
7	PGE	B	523	-	-	0/7/7/7	0/0/0/0
8	P6G	B	524	-	-	0/16/16/16	0/0/0/0
8	P6G	B	525	-	-	0/16/16/16	0/0/0/0
2	ASP	C	501	-	-	0/2/8/8	0/0/0/0
4	DMU	C	505	-	-	0/19/59/59	0/2/2/2
5	1PE	C	506	-	-	0/13/13/13	0/0/0/0
5	1PE	C	507	-	-	0/13/13/13	0/0/0/0
6	PEG	C	508	-	-	0/4/4/4	0/0/0/0
6	PEG	C	509	-	-	0/4/4/4	0/0/0/0
6	PEG	C	510	-	-	0/4/4/4	0/0/0/0
6	PEG	C	511	-	-	0/4/4/4	0/0/0/0
6	PEG	C	512	-	-	0/4/4/4	0/0/0/0
6	PEG	C	513	-	-	0/4/4/4	0/0/0/0
6	PEG	C	514	-	-	0/4/4/4	0/0/0/0
6	PEG	C	515	-	-	0/4/4/4	0/0/0/0
6	PEG	C	516	-	-	0/4/4/4	0/0/0/0
7	PGE	C	517	-	-	0/7/7/7	0/0/0/0
7	PGE	C	518	-	-	0/7/7/7	0/0/0/0
8	P6G	C	519	-	-	0/16/16/16	0/0/0/0
8	P6G	C	520	-	-	0/16/16/16	0/0/0/0

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	505	DMU	C11-C9	-3.29	1.40	1.51
4	B	505	DMU	C11-C9	-3.11	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	505	DMU	C11-C9	-3.10	1.41	1.51
4	C	505	DMU	C5-C7	-2.33	1.46	1.52
4	A	505	DMU	C5-C7	-2.13	1.46	1.52
4	B	505	DMU	C5-C7	-2.12	1.47	1.52
4	C	505	DMU	O3-C5	2.09	1.47	1.43
4	C	505	DMU	O1-C10	2.16	1.47	1.41
4	A	505	DMU	O3-C5	2.17	1.48	1.43
4	C	505	DMU	O5-C4	2.29	1.49	1.44
4	A	505	DMU	C8-C9	2.30	1.57	1.53
4	B	505	DMU	O3-C5	2.31	1.48	1.43
4	C	505	DMU	C8-C9	2.35	1.58	1.53
4	B	505	DMU	O5-C4	2.37	1.50	1.44
4	C	505	DMU	O5-C6	2.40	1.47	1.41
4	A	505	DMU	O1-C10	2.43	1.47	1.41
4	C	505	DMU	O4-C7	2.55	1.48	1.43
4	B	505	DMU	C8-C9	2.56	1.58	1.53
4	A	505	DMU	O4-C7	2.65	1.49	1.43
4	B	505	DMU	O4-C7	2.72	1.49	1.43
4	B	505	DMU	O5-C6	2.73	1.48	1.41
4	A	505	DMU	O5-C4	2.82	1.51	1.44
4	A	505	DMU	O5-C6	3.10	1.49	1.41
4	C	505	DMU	O1-C9	4.63	1.55	1.44
4	B	505	DMU	O1-C9	4.66	1.55	1.44
4	A	505	DMU	O1-C9	4.74	1.55	1.44

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	505	DMU	O55-C2-C1	-3.00	103.82	110.36
4	A	505	DMU	C57-C4-C3	-2.40	106.69	113.24
4	B	505	DMU	C10-O7-C3	-2.18	112.69	118.00
4	B	505	DMU	C57-C4-C3	-2.10	107.52	113.24
4	C	505	DMU	C10-O7-C3	-2.08	112.92	118.00
4	B	505	DMU	C18-O16-C6	2.05	117.38	113.87
4	C	505	DMU	C1-C2-C3	2.33	114.43	109.61
4	A	505	DMU	C6-O5-C4	2.39	118.21	113.72
4	A	505	DMU	C7-C8-C9	2.39	114.43	110.22
4	A	505	DMU	C2-C3-C4	2.48	116.14	110.88
4	B	505	DMU	C7-C8-C9	2.89	115.30	110.22
4	B	505	DMU	O5-C4-C3	3.01	115.91	109.75
4	A	505	DMU	C18-O16-C6	3.16	119.29	113.87
4	C	505	DMU	O5-C4-C3	3.23	116.37	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	505	DMU	O5-C4-C3	3.90	117.73	109.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

21 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	ASP	2	0
5	A	507	1PE	2	0
6	A	511	PEG	1	0
7	A	520	PGE	1	0
7	A	521	PGE	1	0
2	B	501	ASP	3	0
4	B	505	DMU	1	0
5	B	506	1PE	1	0
5	B	507	1PE	1	0
5	B	508	1PE	2	0
6	B	509	PEG	1	0
6	B	522	PEG	2	0
8	B	524	P6G	1	0
8	B	525	P6G	1	0
4	C	505	DMU	2	0
6	C	508	PEG	1	0
6	C	509	PEG	1	0
6	C	510	PEG	2	0
6	C	512	PEG	1	0
6	C	515	PEG	1	0
8	C	520	P6G	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	427/438 (97%)	-0.18	16 (3%) 42 31	55, 84, 153, 243	0
1	B	427/438 (97%)	0.06	27 (6%) 21 13	58, 90, 159, 234	0
1	C	426/438 (97%)	-0.08	27 (6%) 21 13	56, 94, 146, 212	0
All	All	1280/1314 (97%)	-0.07	70 (5%) 26 17	55, 89, 154, 243	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	430	SER	8.7
1	B	127	ALA	8.4
1	B	126	GLU	7.6
1	B	124	ALA	6.7
1	B	11	ASP	5.8
1	B	125	ILE	5.8
1	C	123	LYS	5.7
1	B	16	TRP	5.5
1	C	125	ILE	5.0
1	B	14	VAL	4.9
1	B	123	LYS	4.8
1	C	124	ALA	4.8
1	A	125	ILE	4.7
1	C	16	TRP	4.6
1	B	5	LEU	4.2
1	B	130	PRO	4.1
1	B	15	LEU	3.9
1	A	124	ALA	3.9
1	C	8	ARG	3.9
1	A	375	LEU	3.8
1	B	13	PRO	3.8
1	B	128	GLN	3.7
1	C	122	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	176	GLU	3.6
1	C	422	GLU	3.6
1	C	9	TYR	3.6
1	B	6	LEU	3.5
1	B	129	PRO	3.5
1	A	5	LEU	3.5
1	C	6	LEU	3.4
1	B	431	HIS	3.1
1	B	428	TRP	3.1
1	C	423	LEU	3.1
1	C	378	THR	3.0
1	C	376	ASP	3.0
1	A	121	THR	3.0
1	A	373	VAL	3.0
1	B	421	LYS	2.9
1	A	126	GLU	2.9
1	C	372	SER	2.8
1	C	130	PRO	2.8
1	C	126	GLU	2.8
1	B	44	TYR	2.8
1	A	8	ARG	2.8
1	C	37	TYR	2.8
1	A	6	LEU	2.8
1	C	127	ALA	2.8
1	A	16	TRP	2.7
1	A	340	GLN	2.7
1	C	15	LEU	2.7
1	C	421	LYS	2.6
1	B	430	SER	2.6
1	A	338	LEU	2.6
1	C	10	LEU	2.5
1	A	328	VAL	2.5
1	C	383	VAL	2.5
1	B	342	LEU	2.5
1	B	375	LEU	2.3
1	A	343	VAL	2.2
1	B	131	SER	2.2
1	A	174	ASN	2.2
1	C	5	LEU	2.1
1	B	273	THR	2.1
1	C	11	ASP	2.1
1	B	175	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	403	MET	2.0
1	C	36	GLY	2.0
1	A	273	THR	2.0
1	B	283	THR	2.0
1	C	205	GLN	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	PEG	C	516	7/7	0.83	0.39	22.25	130,134,140,144	0
6	PEG	A	516	7/7	0.90	0.83	20.26	105,108,133,140	0
6	PEG	C	513	7/7	0.67	0.47	17.73	113,118,124,125	0
6	PEG	C	514	7/7	0.62	0.85	14.22	122,124,129,130	0
6	PEG	A	508	7/7	0.93	0.37	13.77	87,89,92,93	0
4	DMU	A	505	33/33	0.73	0.58	13.33	108,156,169,172	0
6	PEG	C	510	7/7	0.88	0.61	13.00	102,109,111,115	0
6	PEG	A	510	7/7	0.81	0.37	9.86	105,106,110,113	0
5	1PE	B	507	16/16	0.63	0.60	9.31	80,121,137,138	0
6	PEG	B	518	7/7	0.89	0.44	9.18	103,114,121,125	0
4	DMU	B	505	33/33	0.85	0.32	7.58	93,142,154,161	0
8	P6G	C	520	19/19	0.76	0.42	6.88	111,120,138,139	0
6	PEG	B	516	7/7	0.78	0.42	6.87	105,118,124,124	0
6	PEG	A	512	7/7	0.84	0.30	6.62	96,112,119,119	0
6	PEG	B	511	7/7	0.76	0.55	5.91	109,122,129,132	0
6	PEG	B	509	7/7	0.89	0.31	5.63	98,101,106,108	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	PEG	B	510	7/7	0.92	0.30	4.94	95,98,100,102	0
8	P6G	C	519	19/19	0.67	0.30	4.64	83,109,120,122	0
7	PGE	A	520	10/10	0.79	0.25	4.38	97,102,114,117	0
5	1PE	A	506	16/16	0.81	0.30	4.34	107,117,124,125	0
5	1PE	B	506	16/16	0.72	0.47	4.25	94,113,121,125	0
7	PGE	A	521	10/10	0.84	0.23	4.23	125,136,140,141	0
8	P6G	B	525	19/19	0.69	0.25	4.13	95,125,130,131	0
4	DMU	C	505	33/33	0.88	0.64	2.94	81,140,149,151	0
5	1PE	C	506	16/16	0.84	0.41	2.49	91,137,155,156	0
8	P6G	A	522	19/19	0.74	0.26	2.42	99,124,135,137	0
7	PGE	C	518	10/10	0.74	0.36	2.11	124,133,135,137	0
6	PEG	B	513	7/7	0.85	0.30	1.96	98,105,121,123	0
8	P6G	B	524	19/19	0.81	0.22	1.74	78,96,112,114	0
7	PGE	A	519	10/10	0.86	0.21	1.68	81,87,108,113	0
3	NA	C	502	1/1	0.99	0.27	1.21	75,75,75,75	0
7	PGE	C	517	10/10	0.86	0.28	0.95	124,132,137,137	0
2	ASP	A	501	9/9	0.95	0.21	0.90	63,70,79,87	0
6	PEG	B	519	7/7	0.69	0.58	0.83	132,133,142,142	0
7	PGE	A	518	10/10	0.90	0.15	0.80	78,94,102,104	0
3	NA	A	502	1/1	0.97	0.24	0.65	63,63,63,63	0
6	PEG	A	513	7/7	0.76	0.26	0.50	125,130,134,137	0
2	ASP	B	501	9/9	0.96	0.19	0.10	63,73,87,104	0
2	ASP	C	501	9/9	0.98	0.18	-0.08	71,76,81,86	0
6	PEG	C	512	7/7	0.81	0.14	-0.08	106,115,124,124	0
3	NA	C	504	1/1	0.88	0.14	-0.26	85,85,85,85	0
6	PEG	B	515	7/7	0.68	0.14	-0.35	109,110,118,120	0
3	NA	A	504	1/1	0.93	0.11	-0.86	86,86,86,86	0
3	NA	B	504	1/1	0.85	0.10	-0.93	112,112,112,112	0
3	NA	A	503	1/1	0.94	0.12	-0.93	65,65,65,65	0
3	NA	B	503	1/1	0.95	0.17	-1.02	61,61,61,61	0
3	NA	C	503	1/1	0.90	0.09	-1.40	66,66,66,66	0
3	NA	B	502	1/1	0.93	0.08	-1.84	68,68,68,68	0
5	1PE	B	508	16/16	0.65	0.36	-	121,139,151,151	0
6	PEG	A	514	7/7	0.63	0.34	-	119,128,132,132	0
6	PEG	C	511	7/7	0.80	0.29	-	115,118,120,120	0
6	PEG	B	522	7/7	0.73	0.56	-	120,125,133,134	0
6	PEG	B	512	7/7	0.79	0.32	-	100,106,116,117	0
6	PEG	B	521	7/7	0.68	0.24	-	124,124,128,129	0
6	PEG	A	511	7/7	0.67	0.21	-	111,116,121,121	0
6	PEG	C	509	7/7	0.78	0.57	-	92,105,108,108	0
6	PEG	C	515	7/7	0.77	0.16	-	121,121,123,123	0
5	1PE	A	507	16/16	0.76	0.30	-	111,124,133,133	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	PEG	B	514	7/7	0.87	0.13	-	104,110,117,118	0
7	PGE	B	523	10/10	0.70	0.23	-	108,130,138,139	0
6	PEG	A	517	7/7	0.83	0.18	-	112,123,133,133	0
6	PEG	C	508	7/7	0.88	0.52	-	93,96,97,97	0
5	1PE	C	507	16/16	0.69	0.15	-	127,143,149,151	0
6	PEG	B	517	7/7	0.73	0.22	-	110,113,118,119	0
6	PEG	A	509	7/7	0.85	0.41	-	87,88,95,95	0
6	PEG	B	520	7/7	0.54	0.30	-	120,125,133,138	0
6	PEG	A	515	7/7	0.80	0.18	-	128,133,136,138	0

6.5 Other polymers [i](#)

There are no such residues in this entry.