



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:55 PM UTC

PDB ID : 6O4D / pdb_00006o4d
Title : Structure of ALDH7A1 mutant W175A complexed with L-pipecolic acid
Authors : Tanner, J.J.; Korasick, D.A.; Laciak, A.R.
Deposited on : 2019-02-28
Resolution : 1.88 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

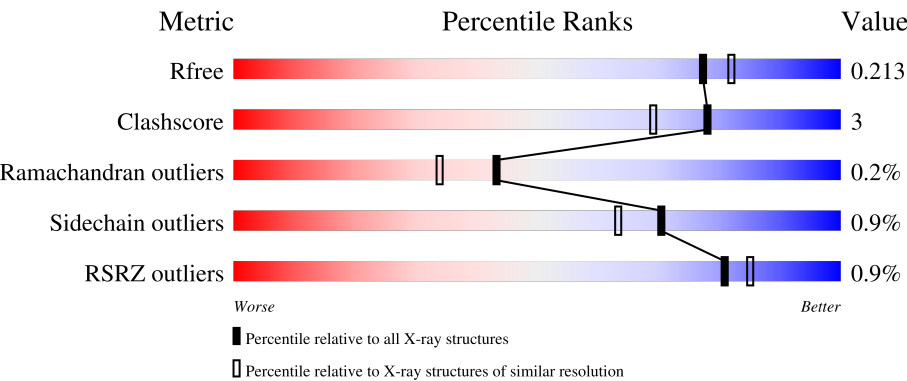
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

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

The reported resolution of this entry is 1.88 Å.

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



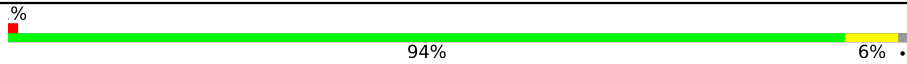
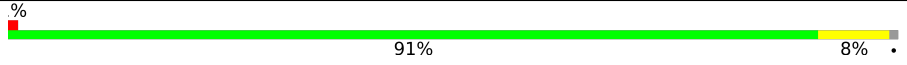
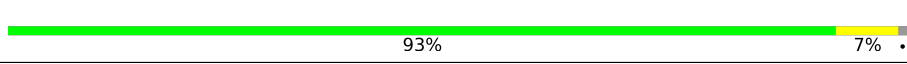
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	513	93%6%.
2	B	513	% 90%9%.
2	G	513	% 91%8%.
3	C	513	2% 91%8%.
3	D	513	% 91%8%.

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Mol	Chain	Length	Quality of chain
3	E	513	 94% 6%
3	F	513	 91% 8%
3	H	513	 93% 7%

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	6PC	A	601	-	X	-	-
4	6PC	B	601	-	X	-	-
4	6PC	C	601	-	X	-	-
4	6PC	D	601	-	X	-	-
4	6PC	E	601	-	X	-	-
4	6PC	F	601	-	X	-	-
4	6PC	G	601	-	X	-	-
4	6PC	H	601	-	X	-	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 33068 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 Alpha-aminoadipic semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	508	Total	C	N	O	S	0	3	0
			3841	2439	666	718	18			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P49419
A	0	HIS	-	expression tag	UNP P49419
A	175	ALA	TRP	engineered mutation	UNP P49419

- Molecule 2 is a protein called Alpha-aminoadipic semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	510	Total	C	N	O	S	0	1	0
			3850	2441	667	724	18			
2	G	508	Total	C	N	O	S	0	2	0
			3855	2445	668	724	18			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP P49419
B	0	HIS	-	expression tag	UNP P49419
B	175	ALA	TRP	engineered mutation	UNP P49419
G	-1	GLY	-	expression tag	UNP P49419
G	0	HIS	-	expression tag	UNP P49419
G	175	ALA	TRP	engineered mutation	UNP P49419

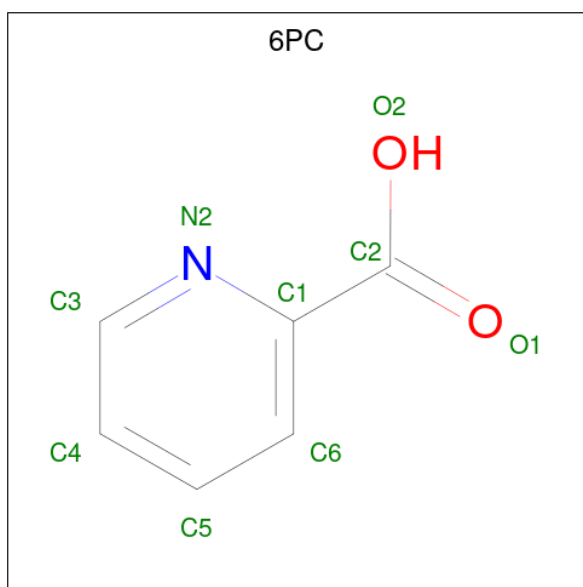
- Molecule 3 is a protein called Alpha-aminoadipic semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	508	Total	C	N	O	S	0	1	0
			3823	2430	662	713	18			
3	D	509	Total	C	N	O	S	0	2	0
			3854	2445	669	722	18			
3	E	509	Total	C	N	O	S	0	1	0
			3843	2441	667	717	18			
3	F	509	Total	C	N	O	S	0	1	0
			3855	2447	669	721	18			
3	H	509	Total	C	N	O	S	0	3	0
			3859	2451	669	721	18			

There are 15 discrepancies between the modelled and reference sequences:

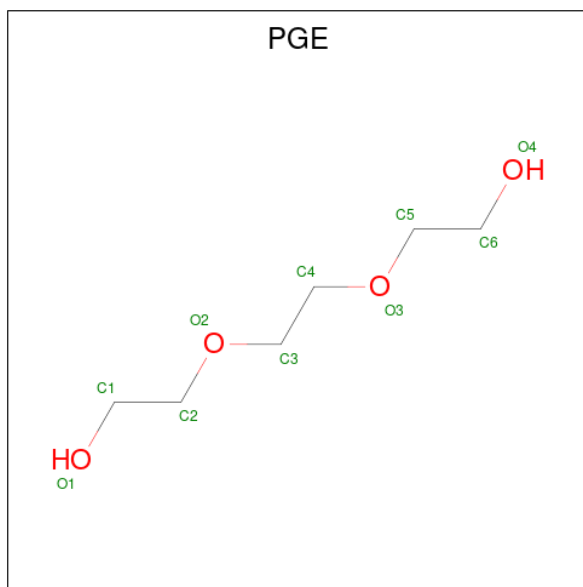
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP P49419
C	0	HIS	-	expression tag	UNP P49419
C	175	ALA	TRP	engineered mutation	UNP P49419
D	-1	GLY	-	expression tag	UNP P49419
D	0	HIS	-	expression tag	UNP P49419
D	175	ALA	TRP	engineered mutation	UNP P49419
E	-1	GLY	-	expression tag	UNP P49419
E	0	HIS	-	expression tag	UNP P49419
E	175	ALA	TRP	engineered mutation	UNP P49419
F	-1	GLY	-	expression tag	UNP P49419
F	0	HIS	-	expression tag	UNP P49419
F	175	ALA	TRP	engineered mutation	UNP P49419
H	-1	GLY	-	expression tag	UNP P49419
H	0	HIS	-	expression tag	UNP P49419
H	175	ALA	TRP	engineered mutation	UNP P49419

- Molecule 4 is PYRIDINE-2-CARBOXYLIC ACID (CCD ID: 6PC) (formula: C₆H₅NO₂).



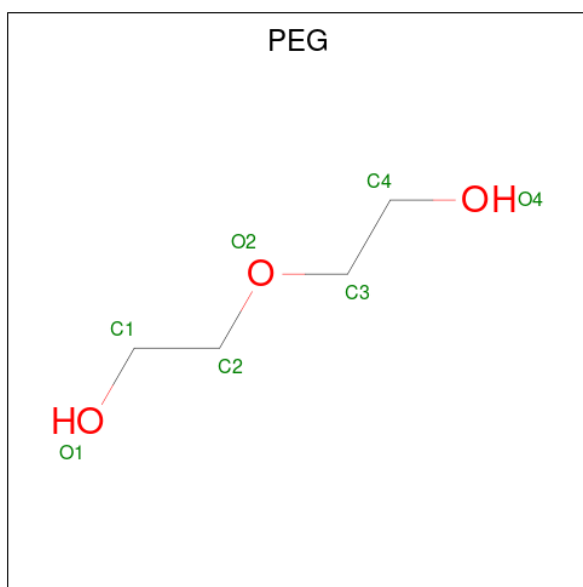
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			9	6	1	2		
4	B	1	Total	C	N	O	0	0
			9	6	1	2		
4	C	1	Total	C	N	O	0	0
			9	6	1	2		
4	D	1	Total	C	N	O	0	0
			9	6	1	2		
4	E	1	Total	C	N	O	0	0
			9	6	1	2		
4	F	1	Total	C	N	O	0	0
			9	6	1	2		
4	G	1	Total	C	N	O	0	0
			9	6	1	2		
4	H	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



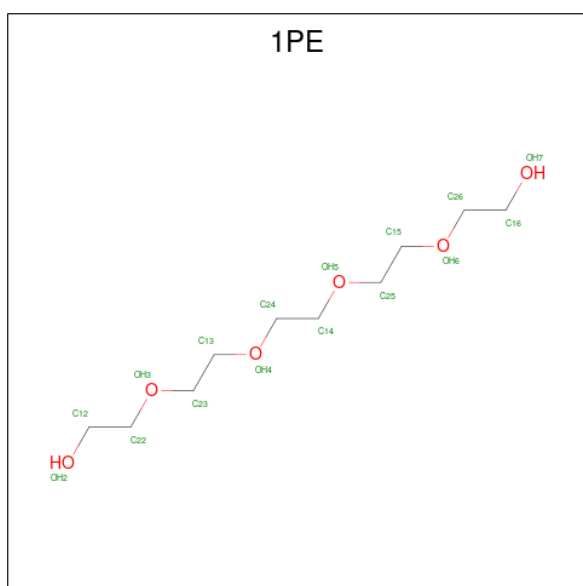
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	H	1	Total	C	O	0	0
			10	6	4		
5	H	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



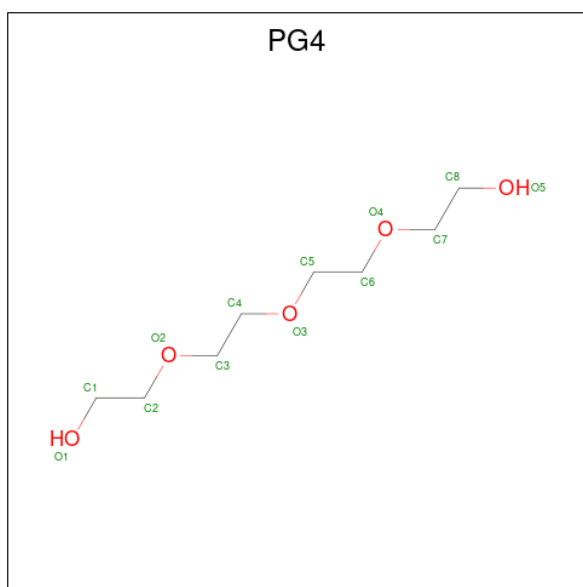
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			7	4	3		
6	F	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



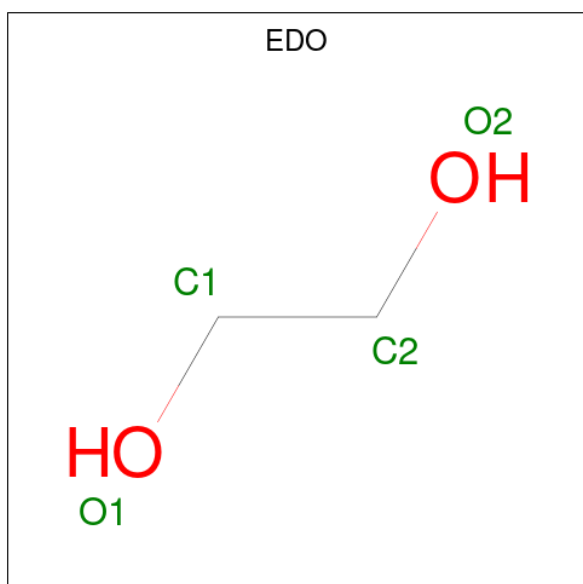
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 8 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	C	O	0	0
			13	8	5		
8	D	1	Total	C	O	0	0
			13	8	5		
8	F	1	Total	C	O	0	0
			13	8	5		
8	G	1	Total	C	O	0	0
			13	8	5		
8	H	1	Total	C	O	0	0
			13	8	5		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	H	1	Total	C	O	0	0
			4	2	2		

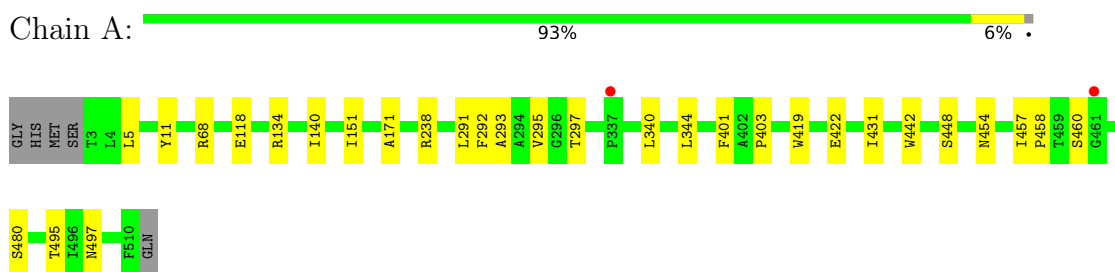
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	261	Total	O	0	0
			261	261		
10	B	263	Total	O	0	0
			263	263		
10	C	241	Total	O	0	1
			242	242		
10	D	223	Total	O	0	0
			223	223		
10	E	273	Total	O	0	0
			273	273		
10	F	254	Total	O	0	0
			254	254		
10	G	276	Total	O	0	1
			277	277		
10	H	264	Total	O	0	0
			264	264		

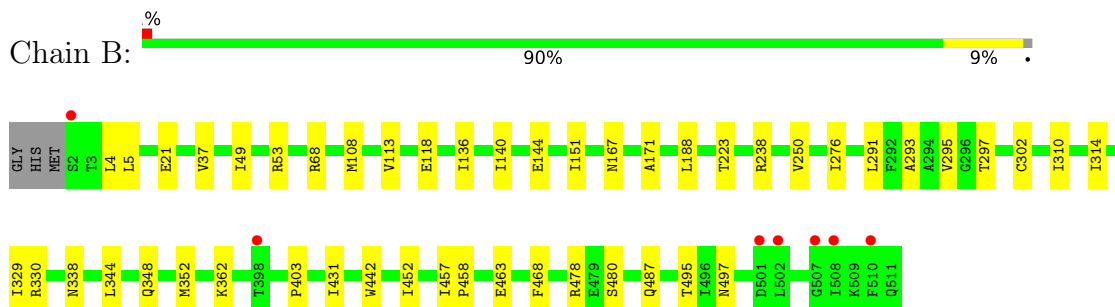
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

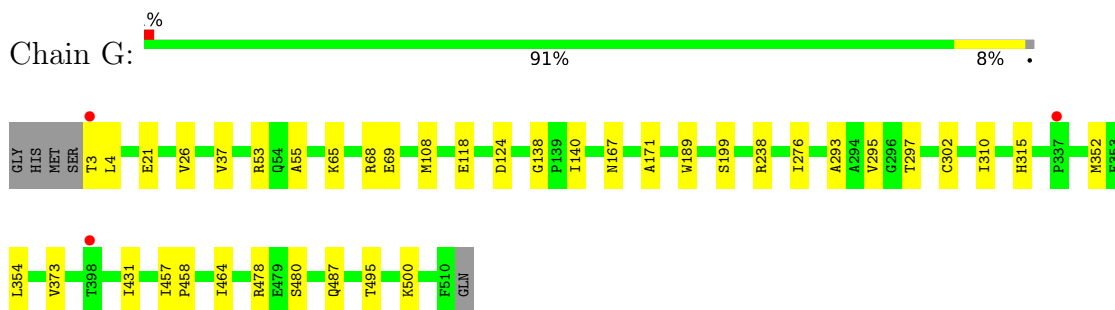
- Molecule 1: Alpha-aminoadipic semialdehyde dehydrogenase



- Molecule 2: Alpha-aminoadipic semialdehyde dehydrogenase

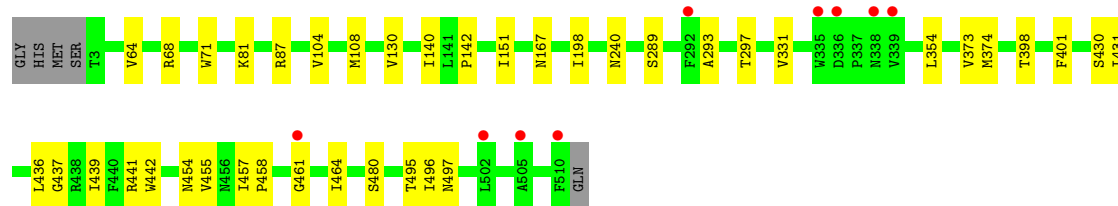


- Molecule 2: Alpha-aminoadipic semialdehyde dehydrogenase

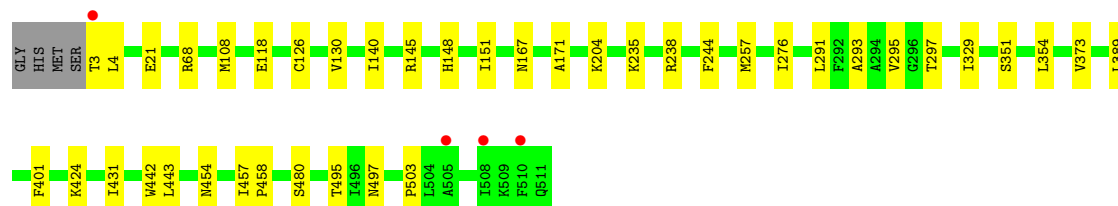
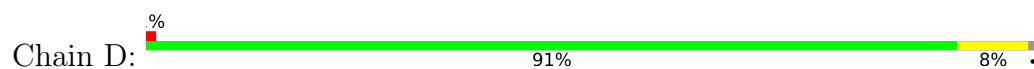


- Molecule 3: Alpha-aminoadipic semialdehyde dehydrogenase

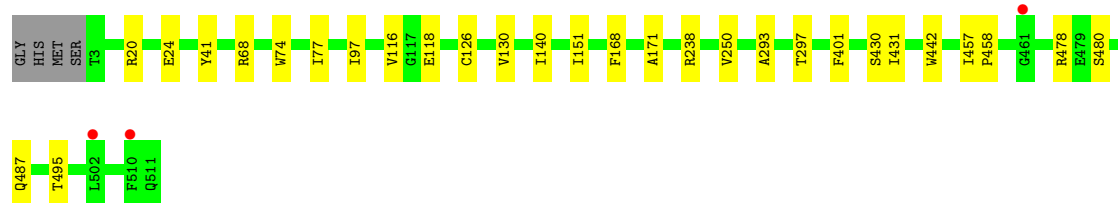
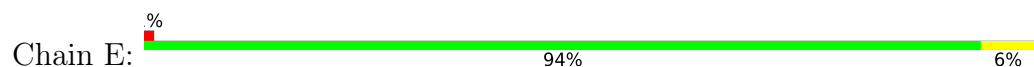




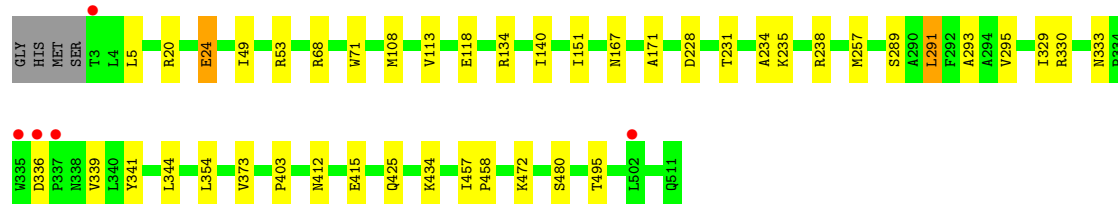
• Molecule 3: Alpha-aminoadipic semialdehyde dehydrogenase



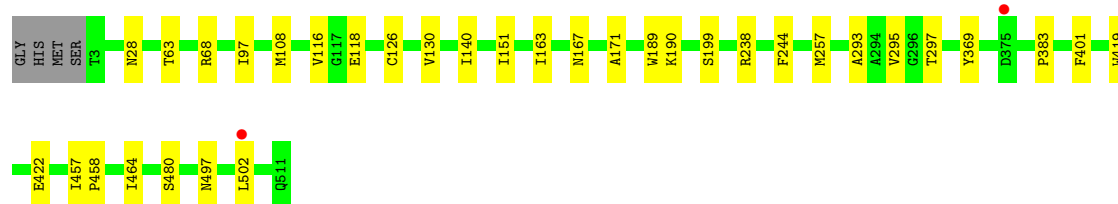
• Molecule 3: Alpha-aminoadipic semialdehyde dehydrogenase



• Molecule 3: Alpha-aminoadipic semialdehyde dehydrogenase



• Molecule 3: Alpha-aminoadipic semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	155.78Å 161.97Å 159.04Å 90.00° 94.88° 90.00°	Depositor
Resolution (Å)	57.96 – 1.88 57.96 – 1.88	Depositor EDS
% Data completeness (in resolution range)	94.7 (57.96-1.88) 90.7 (57.96-1.88)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 1.88Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.172 , 0.215 0.170 , 0.213	Depositor DCC
R_{free} test set	15255 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.501	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	33068	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.30% 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: PG4, 6PC, EDO, OCS, CSO, PEG, 1PE, PGE

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	0.29	0/3919	0.46	0/5318
2	B	0.29	0/3921	0.46	0/5320
2	G	0.28	0/3929	0.45	0/5328
3	C	0.29	0/3904	0.46	0/5301
3	D	0.26	0/3938	0.45	0/5343
3	E	0.29	0/3924	0.47	0/5325
3	F	0.28	0/3936	0.48	0/5339
3	H	0.28	0/3946	0.46	0/5353
All	All	0.28	0/31417	0.46	0/42627

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 [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	3841	0	3812	18	0
2	B	3850	0	3817	31	0
2	G	3855	0	3845	22	0
3	C	3823	0	3788	29	0
3	D	3854	0	3833	28	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	3843	0	3823	18	0
3	F	3855	0	3845	29	0
3	H	3859	0	3849	23	0
4	A	9	0	4	0	0
4	B	9	0	4	0	0
4	C	9	0	4	0	0
4	D	9	0	4	0	0
4	E	9	0	4	1	0
4	F	9	0	4	0	0
4	G	9	0	4	0	0
4	H	9	0	4	0	0
5	A	10	0	14	0	0
5	D	10	0	14	2	0
5	E	20	0	28	2	0
5	H	20	0	28	4	0
6	B	7	0	10	1	0
6	F	7	0	10	0	0
7	B	16	0	22	0	0
8	C	13	0	18	6	0
8	D	13	0	18	2	0
8	F	13	0	18	1	0
8	G	13	0	18	1	0
8	H	13	0	18	2	0
9	H	4	0	6	0	0
10	A	261	0	0	1	0
10	B	263	0	0	2	0
10	C	242	0	0	0	0
10	D	223	0	0	3	0
10	E	273	0	0	0	0
10	F	254	0	0	2	0
10	G	277	0	0	0	0
10	H	264	0	0	0	0
All	All	33068	0	30866	173	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 3.

All (173) 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)
3:C:68:ARG:HH22	8:C:602:PG4:H71	1.31	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:240:ASN:HD21	8:C:602:PG4:H72	1.36	0.89
3:F:293:ALA:HB2	3:F:458:PRO:HB3	1.56	0.87
3:F:235:LYS:HG3	3:F:257:MET:HE3	1.57	0.87
3:E:293:ALA:HB2	3:E:458:PRO:HB3	1.55	0.86
2:B:140:ILE:HD11	3:C:151:ILE:HB	1.65	0.78
3:D:291:LEU:HD11	3:D:329:ILE:HD11	1.70	0.73
1:A:151:ILE:HB	3:D:140:ILE:HD11	1.69	0.72
3:F:291:LEU:HD11	3:F:329:ILE:HD11	1.72	0.71
3:E:140:ILE:HD11	3:H:151:ILE:HB	1.73	0.71
3:F:118:GLU:HG3	3:F:171:ALA:HB2	1.74	0.69
1:A:140:ILE:HD11	3:D:151:ILE:HB	1.74	0.69
3:C:293:ALA:HB2	3:C:458:PRO:HB2	1.75	0.69
3:F:151:ILE:HB	2:G:140:ILE:HD11	1.78	0.64
3:F:234:ALA:HB3	3:F:257:MET:HE2	1.79	0.63
2:G:68:ARG:HB3	8:G:602:PG4:H21	1.82	0.61
2:B:291:LEU:O	2:B:295:VAL:HG22	2.00	0.61
3:H:293:ALA:HB2	3:H:458:PRO:HB3	1.82	0.60
3:H:118:GLU:HG3	3:H:171:ALA:HB2	1.84	0.60
1:A:457:ILE:HD11	2:B:495:THR:HB	1.84	0.60
2:G:37:VAL:HG22	2:G:53:ARG:HG2	1.85	0.59
2:B:293:ALA:HB2	2:B:458:PRO:HB2	1.83	0.59
3:F:336:ASP:O	3:F:339:VAL:HG12	2.03	0.59
3:F:140:ILE:HG22	2:G:138:GLY:O	2.04	0.58
3:D:68:ARG:HB3	8:D:603:PG4:H21	1.84	0.58
3:E:151:ILE:HB	3:H:140:ILE:HD11	1.86	0.57
3:H:383:PRO:HD2	5:H:603:PGE:H52	1.87	0.57
2:B:291:LEU:HD11	2:B:329:ILE:HD11	1.86	0.57
2:G:293:ALA:HB2	2:G:458:PRO:HB2	1.86	0.57
3:C:64:VAL:O	3:C:68:ARG:HG2	2.06	0.56
1:A:68:ARG:HD2	1:A:238:ARG:HB3	1.88	0.55
3:F:20:ARG:NH2	3:F:24:GLU:OE1	2.30	0.55
2:G:478:ARG:O	2:G:487:GLN:NE2	2.34	0.55
3:E:118:GLU:HG3	3:E:171:ALA:HB2	1.89	0.54
3:F:354:LEU:HD21	3:F:373:VAL:HG23	1.90	0.54
3:C:457:ILE:HD11	3:D:495:THR:HB	1.90	0.54
1:A:293:ALA:HB2	1:A:458:PRO:HB2	1.89	0.53
3:D:354:LEU:HD21	3:D:373:VAL:HG23	1.91	0.53
3:H:28:ASN:HB3	3:H:63:THR:HG23	1.89	0.53
3:H:244:PHE:HE1	5:H:605:PGE:H1	1.74	0.53
3:E:74:TRP:HA	3:E:77:ILE:HD12	1.91	0.53
2:B:37:VAL:HG22	2:B:53:ARG:HG2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:354:LEU:HD21	3:C:373:VAL:HG23	1.89	0.52
3:C:495:THR:HB	3:D:457:ILE:HD11	1.92	0.52
2:G:4:LEU:HD12	2:G:21:GLU:HG3	1.92	0.52
3:F:434:LYS:NZ	10:F:703:HOH:O	2.39	0.52
3:D:276:ILE:HB	3:D:431:ILE:HG22	1.92	0.52
3:F:231:THR:HG23	3:F:257:MET:HE1	1.91	0.52
1:A:495:THR:HB	2:B:457:ILE:HD11	1.92	0.51
3:D:293:ALA:HB2	3:D:458:PRO:HB2	1.91	0.51
2:B:151:ILE:HB	3:C:140:ILE:HD11	1.91	0.51
2:B:293:ALA:HB2	2:B:458:PRO:CB	2.40	0.51
2:G:108:MET:HE3	2:G:167:ASN:HA	1.92	0.51
3:D:244:PHE:HE1	5:D:602:PGE:H5	1.75	0.50
3:H:68:ARG:HB3	8:H:604:PG4:H21	1.93	0.50
3:E:457:ILE:HD11	3:F:495:THR:HB	1.93	0.49
2:G:354:LEU:HD21	2:G:373:VAL:HG23	1.94	0.49
3:D:3:THR:N	10:D:711:HOH:O	2.46	0.49
3:F:113:VAL:HG13	10:F:918:HOH:O	2.12	0.49
3:F:339:VAL:HG22	3:F:341:TYR:H	1.78	0.49
1:A:457:ILE:HG23	2:B:497:ASN:HB2	1.94	0.49
3:D:68:ARG:HD2	3:D:238:ARG:HB3	1.94	0.49
3:D:204:LYS:NZ	10:D:709:HOH:O	2.44	0.49
1:A:118:GLU:HG3	1:A:171:ALA:HB2	1.95	0.49
3:C:71:TRP:CE3	8:C:602:PG4:H32	2.48	0.48
1:A:431:ILE:HG23	1:A:442:TRP:CE2	2.49	0.48
2:B:188:LEU:HD11	2:B:223:THR:HG23	1.95	0.48
3:C:454:ASN:HB3	3:C:457:ILE:HG13	1.95	0.48
3:E:20:ARG:NH2	3:E:24:GLU:OE1	2.44	0.48
1:A:5:LEU:HD22	1:A:11:TYR:CE2	2.48	0.48
3:D:235:LYS:HE3	3:D:257:MET:HE2	1.96	0.48
3:H:68:ARG:CZ	8:H:604:PG4:H41	2.44	0.48
2:G:293:ALA:HB2	2:G:458:PRO:CB	2.43	0.48
2:G:495:THR:HB	3:H:457:ILE:HD11	1.95	0.48
3:C:331:VAL:HG21	3:C:374:MET:HE1	1.95	0.48
2:G:276:ILE:HB	2:G:431:ILE:HG22	1.96	0.48
3:F:344:LEU:HD21	3:F:403:PRO:HD3	1.95	0.48
3:H:108:MET:HE3	3:H:167:ASN:HA	1.96	0.47
3:F:333:ASN:O	3:F:339:VAL:HG11	2.13	0.47
3:H:189:TRP:CH2	3:H:199:SER:HA	2.50	0.47
1:A:454:ASN:HB3	1:A:457:ILE:HG13	1.97	0.47
3:H:68:ARG:HD2	3:H:238:ARG:HB3	1.95	0.47
2:B:478:ARG:O	2:B:487:GLN:NE2	2.37	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:87:ARG:HB2	3:C:130:VAL:HG21	1.97	0.47
3:E:495:THR:HB	3:F:457:ILE:HD11	1.96	0.47
3:F:108:MET:HE3	3:F:167:ASN:HA	1.97	0.47
3:F:235:LYS:HE3	3:F:257:MET:HE3	1.96	0.47
2:B:118:GLU:HG3	2:B:171:ALA:HB2	1.95	0.47
3:D:291:LEU:O	3:D:295:VAL:HG22	2.15	0.47
2:B:344:LEU:HD21	2:B:403:PRO:HD3	1.97	0.46
2:G:26:VAL:HG22	2:G:55:ALA:HB2	1.96	0.46
5:D:602:PGE:H62	10:D:878:HOH:O	2.16	0.46
2:B:4:LEU:HD12	2:B:21:GLU:HG3	1.97	0.46
3:C:68:ARG:NH2	8:C:602:PG4:H71	2.14	0.45
3:D:118:GLU:HG3	3:D:171:ALA:HB2	1.98	0.45
2:B:68:ARG:HD2	2:B:238:ARG:HB3	1.98	0.45
3:D:145:ARG:HB3	3:D:503:PRO:HG2	1.98	0.45
1:A:497:ASN:HB2	2:B:457:ILE:HG23	1.98	0.45
3:E:430:SER:OG	3:E:457:ILE:O	2.29	0.45
3:C:430:SER:OG	3:C:457:ILE:O	2.29	0.45
3:H:126:CYS:O	3:H:130:VAL:HG23	2.15	0.45
3:D:108:MET:HE3	3:D:167:ASN:HA	1.99	0.45
3:D:148:HIS:HE1	3:D:503:PRO:HD2	1.81	0.45
2:G:118:GLU:HG3	2:G:171:ALA:HB2	1.98	0.45
3:H:293:ALA:HB2	3:H:458:PRO:CB	2.47	0.44
3:F:235:LYS:HG3	3:F:257:MET:CE	2.38	0.44
3:H:140:ILE:HG23	3:H:151:ILE:HG22	1.99	0.44
2:B:348:GLN:O	2:B:352:MET:HG3	2.17	0.44
2:G:457:ILE:HG23	3:H:497:ASN:HB2	2.00	0.44
3:C:431:ILE:HG23	3:C:442:TRP:CE2	2.52	0.44
3:E:250:VAL:HG13	5:E:603:PGE:H52	2.00	0.44
3:E:168:PHE:CG	4:E:601:6PC:H3	2.53	0.44
3:F:71:TRP:CE3	8:F:602:PG4:H61	2.53	0.44
1:A:344:LEU:HD21	1:A:403:PRO:HD3	1.99	0.43
3:E:68:ARG:HD2	3:E:238:ARG:HB3	2.00	0.43
3:E:431:ILE:HG23	3:E:442:TRP:CE2	2.53	0.43
2:B:310:ILE:HG22	2:B:314:ILE:HG13	2.01	0.43
3:F:68:ARG:HD2	3:F:238:ARG:HB3	2.00	0.43
1:A:134:ARG:HD2	3:C:464:ILE:HD11	2.01	0.43
3:F:228:ASP:OD1	3:F:228:ASP:N	2.51	0.43
2:B:113:VAL:HG13	10:B:898:HOH:O	2.18	0.43
3:C:68:ARG:HH22	8:C:602:PG4:C7	2.17	0.43
2:G:124:ASP:HB3	2:G:464:ILE:HG12	2.01	0.43
2:B:5:LEU:HB2	2:B:49:ILE:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:108:MET:HE3	2:B:167:ASN:HA	2.01	0.43
2:B:250:VAL:HG13	6:B:602:PEG:H22	1.99	0.43
3:F:412:ASN:O	3:F:415:GLU:HG2	2.19	0.42
2:B:452:ILE:HD11	2:B:468:PHE:CD2	2.54	0.42
3:C:439:ILE:HG23	3:C:455:VAL:HG21	1.99	0.42
3:D:454:ASN:HB3	3:D:457:ILE:HG13	2.02	0.42
3:F:134:ARG:HD2	3:H:464:ILE:HD11	2.01	0.42
3:E:126:CYS:O	3:E:130:VAL:HG23	2.18	0.42
1:A:291:LEU:O	1:A:295:VAL:HG22	2.20	0.42
2:B:330:ARG:HD3	2:B:338:ASN:O	2.19	0.42
2:B:431:ILE:HG23	2:B:442:TRP:CE2	2.55	0.42
3:E:97:ILE:HG12	3:E:116:VAL:HG13	2.00	0.42
2:B:362:LYS:HD3	3:E:41:TYR:CZ	2.54	0.42
2:B:463:GLU:HG2	10:B:859:HOH:O	2.20	0.42
3:F:5:LEU:HD12	3:F:49:ILE:HG12	2.02	0.42
2:G:310:ILE:HG21	2:G:315:HIS:HA	2.02	0.42
3:H:369:TYR:OH	5:H:603:PGE:H62	2.19	0.42
3:C:496:ILE:HD11	3:D:443:LEU:HD11	2.01	0.42
3:C:497:ASN:HB2	3:D:457:ILE:HG23	2.02	0.42
8:C:602:PG4:H41	8:C:602:PG4:H62	1.82	0.42
2:G:68:ARG:HD2	2:G:238:ARG:HB3	2.02	0.42
2:G:189:TRP:CH2	2:G:199:SER:HA	2.55	0.42
3:C:108:MET:HE3	3:C:167:ASN:HA	2.02	0.42
3:F:289:SER:HB3	3:F:458:PRO:HG3	2.00	0.41
3:H:163:ILE:HG23	3:H:190:LYS:HE3	2.02	0.41
3:H:257:MET:HE1	5:H:605:PGE:H5	2.00	0.41
1:A:419:TRP:O	1:A:422:GLU:HG2	2.19	0.41
3:H:97:ILE:HG12	3:H:116:VAL:HG13	2.02	0.41
1:A:292:PHE:HB2	10:A:906:HOH:O	2.19	0.41
3:C:457:ILE:HG23	3:D:497:ASN:HB2	2.02	0.41
3:D:4:LEU:HD12	3:D:21:GLU:HG3	2.01	0.41
2:G:65:LYS:O	2:G:69:GLU:HG2	2.19	0.41
2:B:136:ILE:O	3:C:142:PRO:HD3	2.21	0.41
3:C:461:GLY:O	3:D:145:ARG:NH2	2.41	0.41
3:D:68:ARG:CZ	8:D:603:PG4:H31	2.51	0.41
1:A:442:TRP:CH2	1:A:448:SER:HB2	2.56	0.41
2:B:144:GLU:HB2	3:C:81:LYS:HE3	2.02	0.41
2:B:276:ILE:HB	2:B:431:ILE:HG22	2.02	0.41
3:C:104:VAL:HA	3:C:198:ILE:HD11	2.02	0.41
3:C:437:GLY:O	3:C:441:ARG:HG3	2.21	0.41
3:D:126:CYS:O	3:D:130:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:478:ARG:O	3:E:487:GLN:NE2	2.40	0.41
3:F:425:GLN:O	3:F:472:LYS:HE2	2.21	0.41
3:D:431:ILE:HG23	3:D:442:TRP:CE2	2.57	0.40
3:E:68:ARG:CZ	5:E:602:PGE:H6	2.50	0.40
3:H:419:TRP:O	3:H:422:GLU:HG2	2.21	0.40
3:C:289:SER:HB3	3:C:458:PRO:HD3	2.03	0.40
2:G:500:LYS:HD3	2:G:500:LYS:HA	1.82	0.40
2:G:4:LEU:CD1	2:G:21:GLU:HG3	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	508/513 (99%)	493 (97%)	14 (3%)	1 (0%)	43	34
2	B	508/513 (99%)	491 (97%)	16 (3%)	1 (0%)	43	34
2	G	507/513 (99%)	492 (97%)	14 (3%)	1 (0%)	43	34
3	C	507/513 (99%)	490 (97%)	16 (3%)	1 (0%)	43	34
3	D	509/513 (99%)	492 (97%)	16 (3%)	1 (0%)	43	34
3	E	508/513 (99%)	492 (97%)	15 (3%)	1 (0%)	43	34
3	F	508/513 (99%)	493 (97%)	14 (3%)	1 (0%)	43	34
3	H	510/513 (99%)	493 (97%)	16 (3%)	1 (0%)	43	34
All	All	4065/4104 (99%)	3936 (97%)	121 (3%)	8 (0%)	43	34

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	480	SER
1	A	480	SER

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Mol	Chain	Res	Type
3	E	480	SER
3	F	480	SER
2	G	480	SER
3	H	480	SER
2	B	480	SER
3	D	480	SER

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	395/408 (97%)	391 (99%)	4 (1%)	68	60
2	B	397/408 (97%)	396 (100%)	1 (0%)	86	82
2	G	401/408 (98%)	397 (99%)	4 (1%)	68	60
3	C	392/409 (96%)	388 (99%)	4 (1%)	68	60
3	D	400/409 (98%)	395 (99%)	5 (1%)	61	50
3	E	397/409 (97%)	395 (100%)	2 (0%)	81	76
3	F	401/409 (98%)	396 (99%)	5 (1%)	63	53
3	H	401/409 (98%)	397 (99%)	4 (1%)	68	60
All	All	3184/3269 (97%)	3155 (99%)	29 (1%)	70	63

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	297	THR
1	A	340	LEU
1	A	401	PHE
1	A	460	SER
2	B	297	THR
3	C	297	THR
3	C	398	THR
3	C	401	PHE
3	C	436	LEU

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Mol	Chain	Res	Type
3	D	297	THR
3	D	351	SER
3	D	389	LEU
3	D	401	PHE
3	D	424	LYS
3	E	297	THR
3	E	401	PHE
3	F	24	GLU
3	F	53	ARG
3	F	291	LEU
3	F	295	VAL
3	F	330	ARG
2	G	3	THR
2	G	295	VAL
2	G	297	THR
2	G	352	MET
3	H	295	VAL
3	H	297	THR
3	H	401	PHE
3	H	502	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	98	GLN
1	A	120	GLN
1	A	153	GLN
1	A	155	ASN
1	A	177	ASN
1	A	213	ASN
1	A	333	ASN
2	B	46	ASN
2	B	54	GLN
2	B	153	GLN
2	B	155	ASN
2	B	177	ASN
2	B	420	ASN
3	C	155	ASN
3	C	379	ASN
3	C	397	HIS
3	D	46	ASN

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Mol	Chain	Res	Type
3	D	148	HIS
3	D	153	GLN
3	D	155	ASN
3	E	155	ASN
3	E	177	ASN
3	F	155	ASN
3	F	177	ASN
3	F	259	GLN
3	F	397	HIS
2	G	8	GLN
2	G	10	GLN
2	G	46	ASN
2	G	213	ASN
3	H	88	GLN
3	H	120	GLN
3	H	155	ASN
3	H	213	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	OCS	B	302	2	6,8,9	0.94	0	7,11,13	1.65	2 (28%)
2	OCS	G	302	2	6,8,9	0.82	0	7,11,13	1.34	1 (14%)
1	CSO	A	302	1	3,6,7	0.68	0	1,6,8	0.12	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	OCS	B	302	2	-	0/4/7/9	-
2	OCS	G	302	2	-	0/4/7/9	-
1	CSO	A	302	1	-	0/1/5/7	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	302	OCS	OD3-SG-CB	2.79	110.92	106.76
2	B	302	OCS	OD2-SG-CB	2.20	110.23	105.97
2	G	302	OCS	OD2-SG-CB	2.20	110.22	105.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

23 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	$\# Z > 2$
9	EDO	H	602	-	3,3,3	0.50	0	2,2,2	0.26	0
8	PG4	C	602	-	12,12,12	0.51	0	11,11,11	0.35	0
8	PG4	F	602	-	12,12,12	0.53	0	11,11,11	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PGE	H	605	-	9,9,9	0.54	0	8,8,8	0.26	0
8	PG4	G	602	-	12,12,12	0.54	0	11,11,11	0.24	0
5	PGE	D	602	-	9,9,9	0.35	0	8,8,8	0.32	0
4	6PC	A	601	-	9,9,9	5.49	6 (66%)	11,11,11	4.15	8 (72%)
8	PG4	D	603	-	12,12,12	0.52	0	11,11,11	0.24	0
5	PGE	E	603	-	9,9,9	0.29	0	8,8,8	0.41	0
5	PGE	H	603	-	9,9,9	0.35	0	8,8,8	0.27	0
8	PG4	H	604	-	12,12,12	0.57	0	11,11,11	0.48	0
6	PEG	B	602	-	6,6,6	0.50	0	5,5,5	0.24	0
4	6PC	B	601	-	9,9,9	5.52	6 (66%)	11,11,11	4.50	8 (72%)
4	6PC	F	601	-	9,9,9	5.72	6 (66%)	11,11,11	4.24	8 (72%)
4	6PC	H	601	-	9,9,9	5.65	6 (66%)	11,11,11	4.19	9 (81%)
5	PGE	E	602	-	9,9,9	0.29	0	8,8,8	0.24	0
4	6PC	G	601	-	9,9,9	5.51	6 (66%)	11,11,11	4.64	8 (72%)
4	6PC	E	601	-	9,9,9	5.70	7 (77%)	11,11,11	4.39	10 (90%)
4	6PC	C	601	-	9,9,9	5.51	6 (66%)	11,11,11	4.26	8 (72%)
5	PGE	A	602	-	9,9,9	0.36	0	8,8,8	0.33	0
4	6PC	D	601	-	9,9,9	5.53	6 (66%)	11,11,11	4.76	9 (81%)
6	PEG	F	603	-	6,6,6	0.49	0	5,5,5	0.26	0
7	1PE	B	603	-	15,15,15	0.55	0	14,14,14	0.27	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
9	EDO	H	602	-	-	0/1/1/1	-
8	PG4	C	602	-	-	5/10/10/10	-
8	PG4	F	602	-	-	2/10/10/10	-
5	PGE	H	605	-	-	6/7/7/7	-
8	PG4	G	602	-	-	6/10/10/10	-
5	PGE	D	602	-	-	4/7/7/7	-
4	6PC	A	601	-	-	4/4/4/4	0/1/1/1
8	PG4	D	603	-	-	8/10/10/10	-
5	PGE	E	603	-	-	4/7/7/7	-
5	PGE	H	603	-	-	5/7/7/7	-
8	PG4	H	604	-	-	4/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	B	602	-	-	3/4/4/4	-
4	6PC	B	601	-	-	2/4/4/4	0/1/1/1
4	6PC	F	601	-	-	2/4/4/4	0/1/1/1
4	6PC	H	601	-	-	2/4/4/4	0/1/1/1
5	PGE	E	602	-	-	2/7/7/7	-
4	6PC	G	601	-	-	2/4/4/4	0/1/1/1
4	6PC	E	601	-	-	2/4/4/4	0/1/1/1
4	6PC	C	601	-	-	2/4/4/4	0/1/1/1
5	PGE	A	602	-	-	3/7/7/7	-
4	6PC	D	601	-	-	2/4/4/4	0/1/1/1
6	PEG	F	603	-	-	3/4/4/4	-
7	1PE	B	603	-	-	4/13/13/13	-

All (49) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	601	6PC	C1-N2	8.65	1.49	1.34
4	E	601	6PC	C1-N2	8.48	1.49	1.34
4	G	601	6PC	C1-N2	8.46	1.49	1.34
4	H	601	6PC	C1-N2	8.26	1.49	1.34
4	D	601	6PC	C1-N2	8.17	1.49	1.34
4	B	601	6PC	C1-N2	8.09	1.48	1.34
4	A	601	6PC	C1-N2	8.03	1.48	1.34
4	C	601	6PC	C1-N2	8.00	1.48	1.34
4	E	601	6PC	C6-C1	7.88	1.54	1.39
4	F	601	6PC	C6-C1	7.75	1.54	1.39
4	H	601	6PC	C6-C1	7.70	1.53	1.39
4	F	601	6PC	C3-N2	7.41	1.50	1.34
4	C	601	6PC	C6-C1	7.37	1.53	1.39
4	A	601	6PC	C6-C1	7.34	1.53	1.39
4	B	601	6PC	C6-C1	7.34	1.53	1.39
4	D	601	6PC	C6-C1	7.33	1.53	1.39
4	G	601	6PC	C6-C1	7.27	1.53	1.39
4	E	601	6PC	C3-N2	7.24	1.49	1.34
4	D	601	6PC	C3-N2	7.16	1.49	1.34
4	A	601	6PC	C3-N2	7.15	1.49	1.34
4	H	601	6PC	C3-N2	7.15	1.49	1.34
4	B	601	6PC	C3-N2	7.06	1.49	1.34
4	G	601	6PC	C3-N2	6.99	1.49	1.34
4	C	601	6PC	C3-N2	6.97	1.49	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	601	6PC	C5-C6	6.62	1.50	1.38
4	E	601	6PC	C5-C6	6.55	1.50	1.38
4	F	601	6PC	C5-C6	6.51	1.50	1.38
4	D	601	6PC	C5-C6	6.48	1.50	1.38
4	B	601	6PC	C5-C6	6.43	1.49	1.38
4	C	601	6PC	C5-C6	6.41	1.49	1.38
4	A	601	6PC	C5-C6	6.21	1.49	1.38
4	G	601	6PC	C5-C6	6.04	1.49	1.38
4	H	601	6PC	C4-C3	5.71	1.53	1.37
4	A	601	6PC	C4-C3	5.64	1.53	1.37
4	C	601	6PC	C4-C3	5.59	1.53	1.37
4	G	601	6PC	C4-C3	5.59	1.53	1.37
4	B	601	6PC	C4-C3	5.52	1.53	1.37
4	D	601	6PC	C4-C3	5.49	1.53	1.37
4	E	601	6PC	C4-C3	5.48	1.53	1.37
4	F	601	6PC	C4-C3	5.46	1.53	1.37
4	B	601	6PC	C5-C4	5.30	1.49	1.38
4	G	601	6PC	C5-C4	5.21	1.49	1.38
4	C	601	6PC	C5-C4	5.21	1.49	1.38
4	H	601	6PC	C5-C4	5.17	1.49	1.38
4	F	601	6PC	C5-C4	5.10	1.49	1.38
4	D	601	6PC	C5-C4	5.08	1.49	1.38
4	E	601	6PC	C5-C4	5.06	1.49	1.38
4	A	601	6PC	C5-C4	5.01	1.49	1.38
4	E	601	6PC	C1-C2	2.01	1.54	1.50

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	601	6PC	C6-C1-N2	-8.32	111.52	122.74
4	B	601	6PC	C6-C1-N2	-8.09	111.83	122.74
4	D	601	6PC	C6-C1-N2	-7.78	112.25	122.74
4	A	601	6PC	C6-C1-N2	-7.65	112.43	122.74
4	G	601	6PC	C6-C1-N2	-7.55	112.56	122.74
4	G	601	6PC	C5-C6-C1	-7.24	109.98	118.63
4	E	601	6PC	C6-C1-N2	-7.12	113.14	122.74
4	H	601	6PC	C6-C1-N2	-7.06	113.22	122.74
4	F	601	6PC	C6-C1-C2	-6.69	108.66	120.32
4	D	601	6PC	C6-C1-C2	-6.31	109.33	120.32
4	B	601	6PC	C6-C1-C2	-6.27	109.38	120.32
4	F	601	6PC	C6-C1-N2	-6.17	114.42	122.74
4	G	601	6PC	C6-C1-C2	-6.14	109.61	120.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	601	6PC	C4-C3-N2	-6.14	113.70	123.42
4	H	601	6PC	C6-C1-C2	-6.03	109.80	120.32
4	A	601	6PC	C6-C1-C2	-6.02	109.82	120.32
4	E	601	6PC	C4-C3-N2	-5.98	113.96	123.42
4	C	601	6PC	C6-C1-C2	-5.81	110.20	120.32
4	F	601	6PC	C4-C3-N2	-5.76	114.31	123.42
4	E	601	6PC	C4-C5-C6	-5.72	113.18	120.24
4	E	601	6PC	C6-C1-C2	-5.50	110.73	120.32
4	D	601	6PC	C5-C6-C1	-5.49	112.07	118.63
4	D	601	6PC	C4-C5-C6	-5.47	113.49	120.24
4	B	601	6PC	C5-C6-C1	-5.45	112.11	118.63
4	G	601	6PC	C4-C5-C6	-5.43	113.53	120.24
4	B	601	6PC	C4-C3-N2	-5.42	114.85	123.42
4	H	601	6PC	C4-C3-N2	-5.38	114.90	123.42
4	C	601	6PC	C4-C3-N2	-5.36	114.94	123.42
4	G	601	6PC	C4-C3-N2	-5.18	115.22	123.42
4	F	601	6PC	C4-C5-C6	-5.12	113.92	120.24
4	A	601	6PC	C4-C5-C6	-4.96	114.11	120.24
4	H	601	6PC	C5-C6-C1	-4.69	113.02	118.63
4	A	601	6PC	C5-C6-C1	-4.58	113.16	118.63
4	A	601	6PC	C2-C1-N2	-4.50	109.55	117.52
4	C	601	6PC	C5-C6-C1	-4.28	113.51	118.63
4	B	601	6PC	C2-C1-N2	-4.27	109.96	117.52
4	F	601	6PC	C2-C1-N2	-4.22	110.05	117.52
4	B	601	6PC	C4-C5-C6	-4.14	115.13	120.24
4	D	601	6PC	C2-C1-N2	-4.00	110.43	117.52
4	E	601	6PC	C5-C6-C1	-3.94	113.92	118.63
4	C	601	6PC	C4-C5-C6	-3.92	115.40	120.24
4	B	601	6PC	C3-N2-C1	-3.85	111.95	116.93
4	D	601	6PC	C5-C4-C3	-3.82	111.60	118.62
4	E	601	6PC	C5-C4-C3	-3.81	111.62	118.62
4	C	601	6PC	C2-C1-N2	-3.72	110.94	117.52
4	A	601	6PC	C4-C3-N2	-3.71	117.54	123.42
4	H	601	6PC	C2-C1-N2	-3.71	110.94	117.52
4	G	601	6PC	C2-C1-N2	-3.71	110.95	117.52
4	E	601	6PC	C2-C1-N2	-3.71	110.95	117.52
4	C	601	6PC	C3-N2-C1	-3.69	112.17	116.93
4	D	601	6PC	C3-N2-C1	-3.68	112.19	116.93
4	H	601	6PC	C4-C5-C6	-3.67	115.71	120.24
4	F	601	6PC	C5-C4-C3	-3.51	112.17	118.62
4	F	601	6PC	C5-C6-C1	-3.42	114.54	118.63
4	H	601	6PC	C3-N2-C1	-3.08	112.95	116.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	601	6PC	C3-N2-C1	-3.06	112.97	116.93
4	E	601	6PC	C3-N2-C1	-2.88	113.21	116.93
4	H	601	6PC	O2-C2-C1	2.81	121.39	114.71
4	H	601	6PC	C5-C4-C3	-2.69	113.68	118.62
4	F	601	6PC	O2-C2-C1	2.65	121.01	114.71
4	A	601	6PC	O2-C2-C1	2.63	120.94	114.71
4	E	601	6PC	O2-C2-C1	2.36	120.32	114.71
4	G	601	6PC	O2-C2-C1	2.28	120.11	114.71
4	D	601	6PC	O2-C2-C1	2.23	120.01	114.71
4	E	601	6PC	O2-C2-O1	-2.17	118.69	123.35
4	B	601	6PC	O2-C2-C1	2.11	119.72	114.71
4	A	601	6PC	O2-C2-O1	-2.09	118.86	123.35
4	C	601	6PC	O2-C2-C1	2.00	119.47	114.71

There are no chirality outliers.

All (77) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	601	6PC	C6-C1-C2-O2
4	B	601	6PC	C6-C1-C2-O1
4	B	601	6PC	C6-C1-C2-O2
4	C	601	6PC	C6-C1-C2-O1
4	C	601	6PC	C6-C1-C2-O2
4	D	601	6PC	C6-C1-C2-O1
4	D	601	6PC	C6-C1-C2-O2
4	E	601	6PC	C6-C1-C2-O1
4	E	601	6PC	C6-C1-C2-O2
4	F	601	6PC	C6-C1-C2-O1
4	F	601	6PC	C6-C1-C2-O2
4	G	601	6PC	C6-C1-C2-O1
4	G	601	6PC	C6-C1-C2-O2
4	H	601	6PC	C6-C1-C2-O1
4	H	601	6PC	C6-C1-C2-O2
5	D	602	PGE	C1-C2-O2-C3
8	C	602	PG4	C4-C3-O2-C2
5	E	603	PGE	O2-C3-C4-O3
5	H	603	PGE	O2-C3-C4-O3
4	A	601	6PC	C6-C1-C2-O1
8	H	604	PG4	O4-C7-C8-O5
8	C	602	PG4	C6-C5-O3-C4
7	B	603	1PE	OH5-C14-C24-OH4
6	B	602	PEG	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
8	D	603	PG4	O2-C3-C4-O3
8	H	604	PG4	O3-C5-C6-O4
5	A	602	PGE	O1-C1-C2-O2
5	H	605	PGE	O2-C3-C4-O3
7	B	603	1PE	OH7-C16-C26-OH6
8	C	602	PG4	O3-C5-C6-O4
5	D	602	PGE	O2-C3-C4-O3
8	D	603	PG4	O4-C7-C8-O5
8	G	602	PG4	O1-C1-C2-O2
5	E	602	PGE	O3-C5-C6-O4
6	B	602	PEG	O2-C3-C4-O4
6	F	603	PEG	O2-C3-C4-O4
8	G	602	PG4	O4-C7-C8-O5
8	D	603	PG4	O3-C5-C6-O4
8	C	602	PG4	O1-C1-C2-O2
5	H	603	PGE	C6-C5-O3-C4
5	E	603	PGE	C3-C4-O3-C5
6	F	603	PEG	C1-C2-O2-C3
8	D	603	PG4	C8-C7-O4-C6
8	H	604	PG4	C6-C5-O3-C4
5	H	605	PGE	C6-C5-O3-C4
5	H	603	PGE	O1-C1-C2-O2
5	H	605	PGE	O3-C5-C6-O4
6	F	603	PEG	O1-C1-C2-O2
5	H	605	PGE	C1-C2-O2-C3
5	H	603	PGE	C4-C3-O2-C2
5	D	602	PGE	C3-C4-O3-C5
8	F	602	PG4	C1-C2-O2-C3
6	B	602	PEG	C1-C2-O2-C3
5	E	602	PGE	O1-C1-C2-O2
7	B	603	1PE	C24-C14-OH5-C25
5	H	603	PGE	C3-C4-O3-C5
8	F	602	PG4	C4-C3-O2-C2
8	H	604	PG4	C4-C3-O2-C2
7	B	603	1PE	C25-C15-OH6-C26
8	D	603	PG4	C3-C4-O3-C5
5	E	603	PGE	C4-C3-O2-C2
5	A	602	PGE	O3-C5-C6-O4
8	D	603	PG4	C1-C2-O2-C3
8	G	602	PG4	C4-C3-O2-C2
8	D	603	PG4	O1-C1-C2-O2
4	A	601	6PC	N2-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	H	605	PGE	C3-C4-O3-C5
5	H	605	PGE	C4-C3-O2-C2
8	G	602	PG4	O2-C3-C4-O3
8	D	603	PG4	C6-C5-O3-C4
5	A	602	PGE	C4-C3-O2-C2
8	G	602	PG4	C6-C5-O3-C4
4	A	601	6PC	N2-C1-C2-O1
5	E	603	PGE	O1-C1-C2-O2
5	D	602	PGE	O3-C5-C6-O4
8	C	602	PG4	O2-C3-C4-O3
8	G	602	PG4	C1-C2-O2-C3

There are no ring outliers.

12 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	602	PG4	6	0
8	F	602	PG4	1	0
5	H	605	PGE	2	0
8	G	602	PG4	1	0
5	D	602	PGE	2	0
8	D	603	PG4	2	0
5	E	603	PGE	1	0
5	H	603	PGE	2	0
8	H	604	PG4	2	0
6	B	602	PEG	1	0
5	E	602	PGE	1	0
4	E	601	6PC	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(Å²)	Q<0.9	
1	A	507/513 (98%)	-0.26	2 (0%)	88	92	9, 26, 42, 62	3 (0%)
2	B	509/513 (99%)	-0.25	7 (1%)	73	79	15, 26, 43, 75	1 (0%)
2	G	507/513 (98%)	-0.29	3 (0%)	85	89	15, 25, 41, 64	2 (0%)
3	C	508/513 (99%)	-0.22	9 (1%)	67	73	11, 25, 48, 69	1 (0%)
3	D	509/513 (99%)	-0.10	4 (0%)	82	86	12, 30, 54, 77	2 (0%)
3	E	509/513 (99%)	-0.29	3 (0%)	85	89	14, 26, 40, 61	1 (0%)
3	F	509/513 (99%)	-0.19	5 (0%)	79	83	14, 27, 44, 64	1 (0%)
3	H	509/513 (99%)	-0.29	2 (0%)	88	92	11, 26, 43, 57	3 (0%)
All	All	4067/4104 (99%)	-0.24	35 (0%)	81	85	9, 26, 45, 77	14 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	335	TRP	5.0
3	D	510	PHE	4.5
3	D	508	ILE	4.4
2	B	510	PHE	3.9
3	E	510	PHE	3.5
3	C	502	LEU	3.2
3	C	510	PHE	3.2
3	C	338	ASN	3.0
3	F	502	LEU	3.0
3	C	292	PHE	2.8
3	C	461	GLY	2.8
2	B	2	SER	2.7
2	G	3	THR	2.7
2	B	502	LEU	2.6
3	F	336	ASP	2.6
2	B	501	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	398	THR	2.5
3	F	335	TRP	2.5
3	C	336	ASP	2.4
3	D	505	ALA	2.4
1	A	461	GLY	2.4
3	E	502	LEU	2.4
1	A	337	PRO	2.3
3	F	337	PRO	2.3
2	B	508	ILE	2.3
3	C	339	VAL	2.3
3	C	505	ALA	2.2
3	E	461	GLY	2.2
3	H	502	LEU	2.2
2	G	337	PRO	2.2
3	H	375	ASP	2.2
2	G	398	THR	2.2
2	B	507	GLY	2.1
3	D	3	THR	2.0
3	F	3	THR	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
1	CSO	A	302	7/8	0.94	0.07	23,26,40,40	0
2	OCS	G	302	9/10	0.96	0.09	18,28,41,53	0
2	OCS	B	302	9/10	0.97	0.10	17,20,44,51	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	PGE	D	602	10/10	0.82	0.16	51,54,55,55	0
5	PGE	E	603	10/10	0.85	0.13	39,46,50,53	0
5	PGE	H	603	10/10	0.86	0.15	43,48,51,53	0
4	6PC	F	601	9/9	0.88	0.11	41,44,47,48	0
5	PGE	H	605	10/10	0.88	0.13	42,48,54,58	0
8	PG4	H	604	13/13	0.88	0.12	22,39,43,46	0
4	6PC	E	601	9/9	0.89	0.11	39,41,43,44	0
5	PGE	E	602	10/10	0.89	0.13	31,40,46,50	0
4	6PC	A	601	9/9	0.89	0.11	38,43,53,55	0
4	6PC	C	601	9/9	0.90	0.12	32,35,44,45	0
8	PG4	C	602	13/13	0.90	0.10	30,36,44,46	0
4	6PC	D	601	9/9	0.90	0.11	41,45,49,51	0
8	PG4	G	602	13/13	0.91	0.11	36,44,59,60	0
4	6PC	B	601	9/9	0.91	0.10	33,36,40,43	0
9	EDO	H	602	4/4	0.91	0.15	40,43,43,44	0
8	PG4	F	602	13/13	0.92	0.11	36,43,51,51	0
7	1PE	B	603	16/16	0.92	0.11	36,42,58,60	0
6	PEG	F	603	7/7	0.92	0.12	46,47,54,56	0
8	PG4	D	603	13/13	0.92	0.10	37,42,46,48	0
6	PEG	B	602	7/7	0.93	0.10	37,46,50,50	0
5	PGE	A	602	10/10	0.93	0.10	34,38,44,46	0
4	6PC	G	601	9/9	0.93	0.10	25,32,35,35	0
4	6PC	H	601	9/9	0.94	0.08	24,29,33,34	0

6.5 Other polymers ⓘ

There are no such residues in this entry.