



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 18, 2025 – 04:07 pm BST

PDB ID : 9FOI / pdb_00009foi
Title : Structure of human KCTD1
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Deposited on : 2024-06-11
Resolution : 2.71 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 Phenix2.0rc1
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

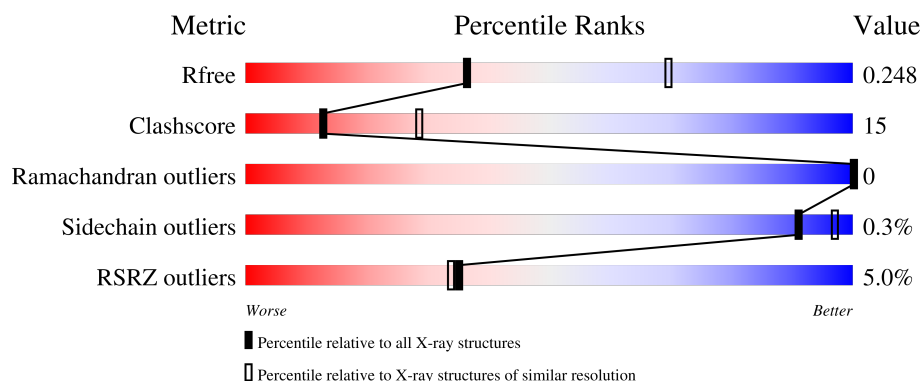
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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}	164625	4050 (2.74-2.70)
Clashscore	180529	4439 (2.74-2.70)
Ramachandran outliers	177936	4374 (2.74-2.70)
Sidechain outliers	177891	4375 (2.74-2.70)
RSRZ outliers	164620	4050 (2.74-2.70)

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	258	<div> <div>5%</div> <div> <div>61%</div> <div>21%</div> <div>18%</div> </div> </div>
1	B	258	<div> <div>4%</div> <div> <div>57%</div> <div>24%</div> <div>19%</div> </div> </div>
1	C	258	<div> <div>4%</div> <div> <div>57%</div> <div>26%</div> <div>18%</div> </div> </div>
1	D	258	<div> <div>3%</div> <div> <div>55%</div> <div>26%</div> <div>20%</div> </div> </div>
1	E	258	<div> <div>5%</div> <div> <div>58%</div> <div>22%</div> <div>19%</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	SIN	C	302	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8500 atoms, of which 32 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 BTB/POZ domain-containing protein KCTD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	2	0
			1701	1082	292	319	8			
1	B	208	Total	C	N	O	S	0	0	0
			1653	1056	282	307	8			
1	C	212	Total	C	N	O	S	0	0	0
			1679	1069	289	313	8			
1	D	207	Total	C	N	O	S	0	2	0
			1655	1053	285	309	8			
1	E	209	Total	C	N	O	S	0	0	0
			1664	1058	286	312	8			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q719H9
A	64	ASP	GLU	conflict	UNP Q719H9
B	0	SER	-	expression tag	UNP Q719H9
B	64	ASP	GLU	conflict	UNP Q719H9
C	0	SER	-	expression tag	UNP Q719H9
C	64	ASP	GLU	conflict	UNP Q719H9
D	0	SER	-	expression tag	UNP Q719H9
D	64	ASP	GLU	conflict	UNP Q719H9
E	0	SER	-	expression tag	UNP Q719H9
E	64	ASP	GLU	conflict	UNP Q719H9

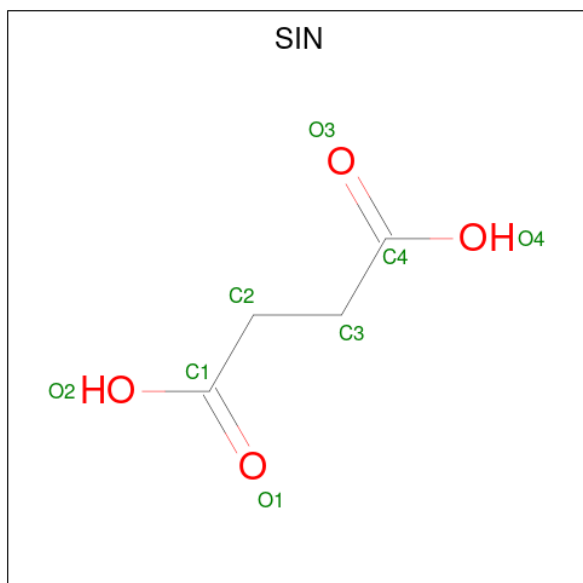
- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	I	0	0
			1	1		
2	C	1	Total	I	0	0
			1	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Na	0	0
			2	2		
3	D	1	Total	Na	0	0
			1	1		

- Molecule 4 is SUCCINIC ACID (CCD ID: SIN) (formula: C₄H₆O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			12	4	4	4		
4	B	1	Total	C	H	O	0	0
			12	4	4	4		
4	C	1	Total	C	H	O	0	0
			12	4	4	4		
4	D	1	Total	C	H	O	0	0
			12	4	4	4		
4	E	1	Total	C	H	O	0	0
			12	4	4	4		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	H	O	0	0
			10	2	6	2		
5	E	1	Total	C	H	O	0	0
			10	2	6	2		

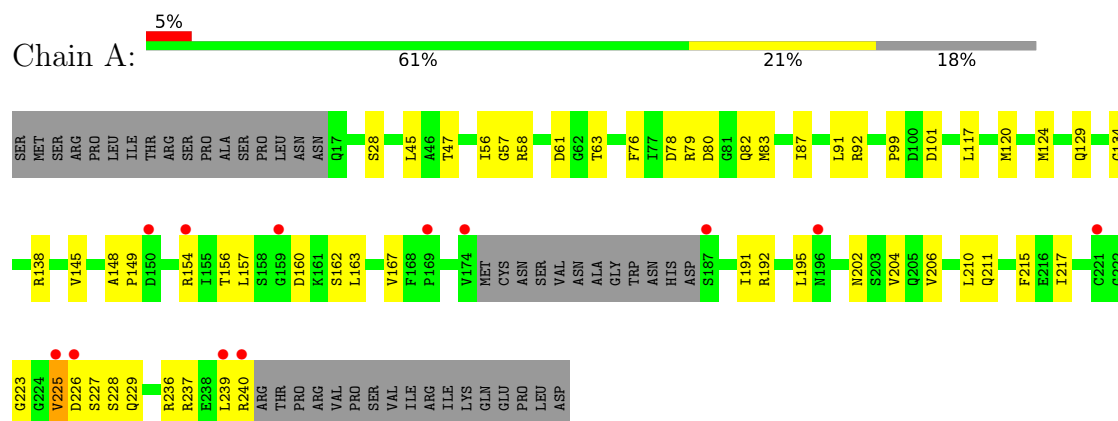
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	14	Total	O	0	0
			14	14		
6	B	8	Total	O	0	0
			8	8		
6	C	9	Total	O	0	0
			9	9		
6	D	11	Total	O	0	0
			11	11		
6	E	21	Total	O	0	0
			21	21		

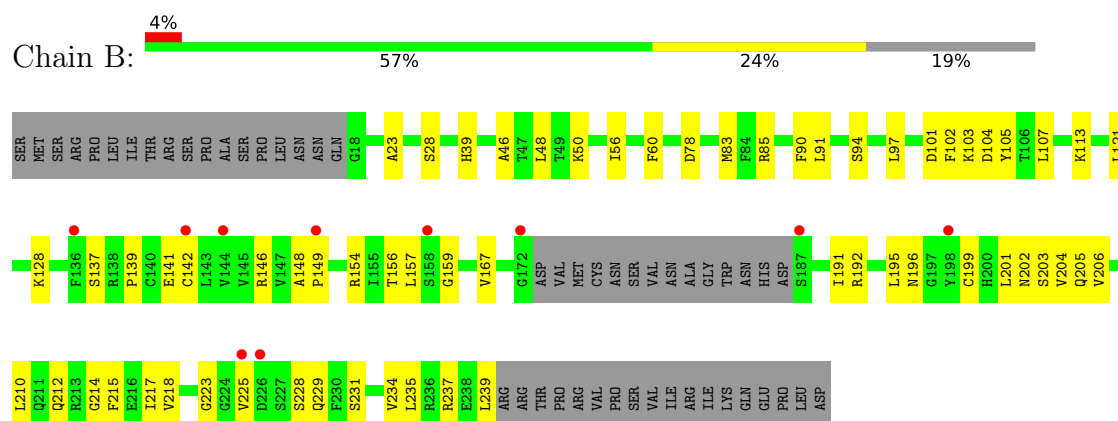
3 Residue-property plots [i](#)

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.

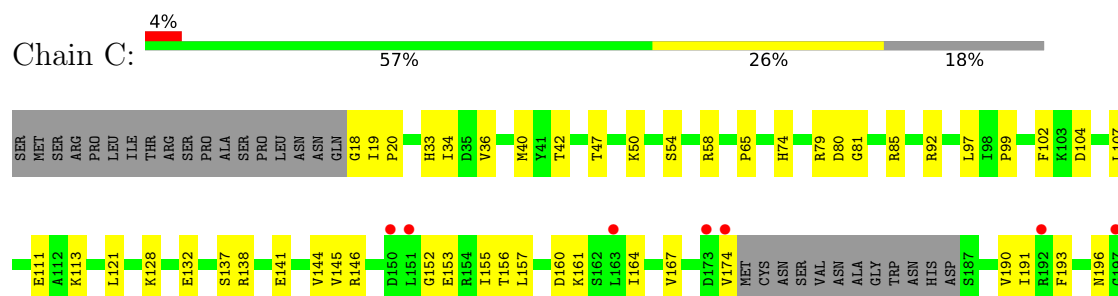
- Molecule 1: BTB/POZ domain-containing protein KCTD1

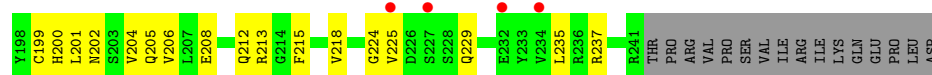


- Molecule 1: BTB/POZ domain-containing protein KCTD1

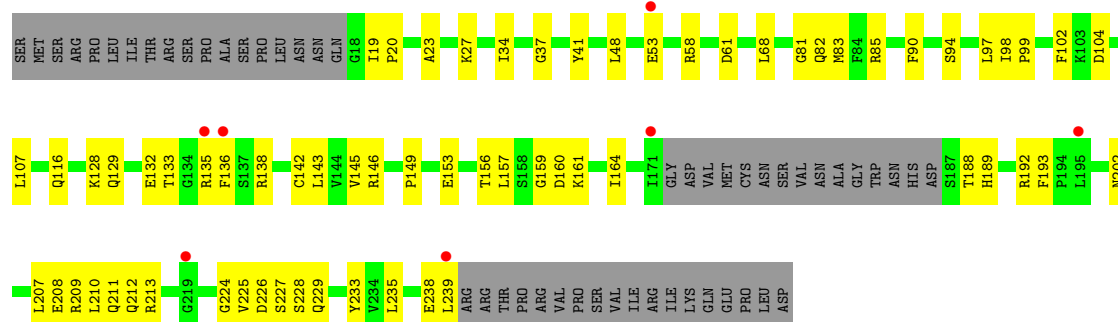


- Molecule 1: BTB/POZ domain-containing protein KCTD1

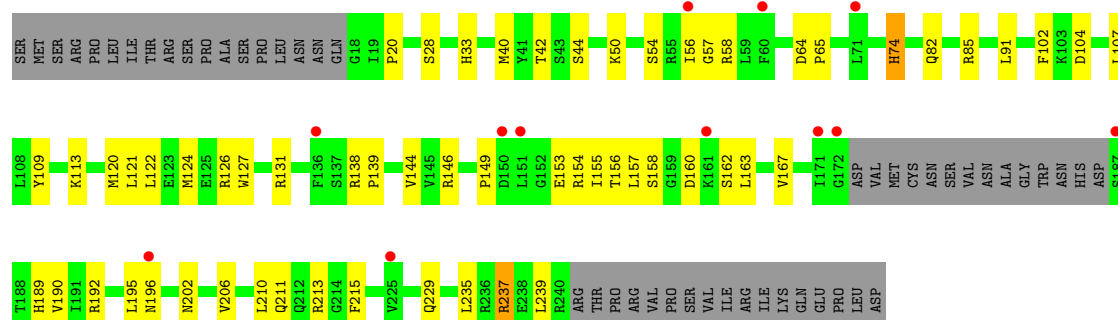




- Molecule 1: BTB/POZ domain-containing protein KCTD1



- Molecule 1: BTB/POZ domain-containing protein KCTD1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.56Å 96.63Å 116.00Å 90.00° 98.09° 90.00°	Depositor
Resolution (Å)	66.89 – 2.71 66.89 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (66.89-2.71) 99.8 (66.89-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.207 , 0.247 0.206 , 0.248	Depositor DCC
R_{free} test set	1949 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	57.0	Xtriage
Anisotropy	0.380	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8500	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.59% 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: EDO, SIN, NA, IOD

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	0.23	0/1743	0.46	0/2357
1	B	0.21	0/1689	0.45	0/2285
1	C	0.22	0/1715	0.44	0/2320
1	D	0.23	0/1697	0.44	0/2296
1	E	0.22	0/1700	0.44	0/2299
All	All	0.22	0/8544	0.45	0/11557

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	1701	0	1666	50	0
1	B	1653	0	1624	79	0
1	C	1679	0	1636	56	0
1	D	1655	0	1609	56	0
1	E	1664	0	1623	49	0
2	A	1	0	0	0	0
2	C	1	0	0	0	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	1	0	0	0	0
4	A	8	4	4	1	0
4	B	8	4	4	1	0
4	C	8	4	4	5	0
4	D	8	4	4	0	0
4	E	8	4	4	2	0
5	B	4	6	6	0	0
5	E	4	6	6	2	0
6	A	14	0	0	1	0
6	B	8	0	0	0	0
6	C	9	0	0	0	0
6	D	11	0	0	0	0
6	E	21	0	0	0	0
All	All	8468	32	8190	247	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 15.

All (247) 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:157:LEU:HD11	1:A:195:LEU:HD22	1.54	0.86
1:B:141:GLU:HG3	1:B:239:LEU:HD23	1.61	0.82
1:A:225:VAL:HG12	1:B:225:VAL:HG13	1.62	0.82
1:B:85:ARG:HG3	4:C:302:SIN:H32	1.62	0.79
1:A:149:PRO:HB3	1:B:192:ARG:HH12	1.48	0.77
1:B:39:HIS:HB2	4:B:301:SIN:H32	1.68	0.76
1:B:156:THR:HB	1:B:192:ARG:HA	1.67	0.76
1:A:202:ASN:O	1:A:206:VAL:HG23	1.87	0.75
1:C:80:ASP:HA	4:C:302:SIN:O1	1.86	0.74
1:A:61:ASP:OD1	1:A:63:THR:HG23	1.87	0.73
1:E:146:ARG:HB2	1:E:156:THR:OG1	1.88	0.73
1:D:188:THR:HG23	1:D:189:HIS:ND1	2.06	0.71
1:B:141:GLU:CG	1:B:239:LEU:HD23	2.20	0.70
1:E:167:VAL:HG11	1:E:210:LEU:HD23	1.73	0.69
1:B:225:VAL:O	1:B:229:GLN:HG3	1.91	0.69
1:C:99:PRO:HD2	1:C:102:PHE:CD1	2.27	0.69
1:A:149:PRO:HG3	1:B:192:ARG:HH22	1.57	0.69
1:A:145:VAL:HG22	1:A:157:LEU:CD2	2.23	0.68
1:D:85[B]:ARG:HE	1:E:82:GLN:HG3	1.57	0.68
1:A:149:PRO:HB3	1:B:192:ARG:NH1	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:GLY:O	1:B:229:GLN:HA	1.94	0.67
1:B:217:ILE:HG12	1:B:235:LEU:CD2	2.25	0.67
1:D:149:PRO:HD2	1:D:229:GLN:O	1.95	0.65
1:C:204:VAL:HG11	1:D:159:GLY:HA2	1.79	0.65
1:B:141:GLU:CG	1:B:239:LEU:CD2	2.75	0.65
1:D:58:ARG:NH2	1:D:116:GLN:OE1	2.30	0.64
1:D:34:ILE:HD11	1:D:48:LEU:HD11	1.78	0.64
1:A:58:ARG:HB3	1:A:63:THR:OG1	1.98	0.64
1:B:225:VAL:HG22	1:B:228:SER:HB3	1.80	0.64
1:C:225:VAL:C	1:C:229:GLN:HG3	2.23	0.63
1:A:225:VAL:HG22	1:A:228:SER:OG	1.98	0.63
1:C:145:VAL:HG22	1:C:157:LEU:CD1	2.29	0.63
1:D:212:GLN:HA	1:E:138:ARG:O	1.98	0.63
1:D:97:LEU:HB3	1:E:107:LEU:HG	1.81	0.63
1:B:157:LEU:HD22	1:B:195:LEU:HD23	1.80	0.63
1:B:234:VAL:C	1:B:235:LEU:HD23	2.23	0.62
1:C:146:ARG:HB2	1:C:156:THR:OG1	2.00	0.62
1:B:154:ARG:HA	1:B:196:ASN:ND2	2.15	0.61
1:B:83:MET:HG3	1:B:107:LEU:HG	1.81	0.61
1:C:202:ASN:O	1:C:206:VAL:HG23	1.99	0.61
6:A:403:HOH:O	1:E:20:PRO:HA	2.01	0.60
1:B:156:THR:HB	1:B:192:ARG:CA	2.30	0.60
1:A:217:ILE:HD12	1:B:218:VAL:HG23	1.83	0.60
1:D:132:GLU:O	1:D:135:ARG:HG2	2.00	0.60
1:C:167:VAL:HG13	1:C:213:ARG:HG3	1.83	0.60
1:A:239:LEU:HD12	1:A:240:ARG:O	2.01	0.60
1:A:149:PRO:HG2	1:A:229:GLN:CG	2.32	0.59
1:E:154:ARG:HA	1:E:196:ASN:HD21	1.65	0.59
1:B:202:ASN:O	1:B:206:VAL:HG23	2.02	0.59
1:E:153:GLU:O	1:E:155:ILE:HG13	2.02	0.59
1:C:190:VAL:HG21	1:C:193:PHE:HB2	1.84	0.59
1:B:214:GLY:O	1:B:237:ARG:HG3	2.03	0.59
1:D:157:LEU:HD22	1:D:164:ILE:CD1	2.32	0.59
1:A:149:PRO:HG2	1:A:229:GLN:HG3	1.85	0.59
1:C:33:HIS:HB2	1:C:74:HIS:HB3	1.85	0.59
1:B:83:MET:HA	1:B:83:MET:HE2	1.85	0.58
1:B:210:LEU:O	1:B:215:PHE:HB2	2.04	0.58
1:A:76:PHE:CE1	1:E:42:THR:HG22	2.39	0.58
1:C:225:VAL:O	1:C:229:GLN:HG3	2.04	0.57
1:B:217:ILE:HG12	1:B:235:LEU:HD21	1.85	0.57
1:A:148:ALA:HB3	1:A:154:ARG:HB2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLY:O	1:A:229:GLN:HA	2.05	0.57
1:D:129:GLN:O	1:D:133:THR:HG23	2.05	0.57
1:D:156:THR:HA	1:D:193:PHE:O	2.05	0.57
1:A:160:ASP:OD1	1:A:162:SER:HB3	2.04	0.57
1:A:157:LEU:CD1	1:A:195:LEU:HD22	2.29	0.57
1:B:102:PHE:CZ	1:B:104:ASP:HB2	2.40	0.56
1:B:212:GLN:HA	1:C:138:ARG:O	2.05	0.56
1:B:202:ASN:ND2	1:B:204:VAL:HG22	2.20	0.56
1:C:19:ILE:HD11	1:D:27:LYS:HE2	1.86	0.56
1:C:33:HIS:O	1:C:74:HIS:HB2	2.06	0.56
1:D:157:LEU:HD22	1:D:164:ILE:HG12	1.88	0.56
1:B:141:GLU:CD	1:B:239:LEU:CD2	2.78	0.55
1:A:82:GLN:HG3	4:A:304:SIN:H21	1.89	0.55
1:B:137:SER:O	1:B:139:PRO:HD3	2.07	0.55
1:C:79:ARG:HD2	1:C:111:GLU:OE1	2.08	0.54
1:A:120:MET:O	1:A:124:MET:HG3	2.06	0.54
1:D:99:PRO:HG3	1:E:104:ASP:OD1	2.06	0.54
1:B:101:ASP:O	1:B:103:LYS:HG3	2.08	0.54
1:C:128:LYS:O	1:C:132:GLU:HB2	2.08	0.54
1:E:144:VAL:HB	1:E:158:SER:HB2	1.90	0.54
1:B:146:ARG:HA	1:B:231:SER:O	2.07	0.54
1:D:210:LEU:HD13	1:D:235:LEU:HD22	1.89	0.53
1:E:215:PHE:CD1	1:E:237:ARG:HB2	2.43	0.53
1:C:102:PHE:CZ	1:C:104:ASP:HB2	2.43	0.53
1:C:141:GLU:OE2	1:C:237:ARG:HD2	2.09	0.53
1:E:158:SER:HA	1:E:190:VAL:O	2.07	0.53
1:E:40:MET:HG3	5:E:302:EDO:H11	1.90	0.53
1:D:142:CYS:O	1:D:143:LEU:HD23	2.08	0.53
1:D:208:GLU:O	1:D:212:GLN:HG3	2.09	0.53
1:B:215:PHE:CD2	1:B:237:ARG:HB2	2.44	0.53
1:B:97:LEU:CB	1:C:107:LEU:HD13	2.40	0.52
1:B:239:LEU:C	1:B:239:LEU:HD12	2.34	0.52
1:A:99:PRO:HG3	1:B:104:ASP:OD1	2.10	0.52
1:E:157:LEU:HD12	1:E:157:LEU:N	2.23	0.52
1:E:113:LYS:HD2	1:E:121:LEU:HD11	1.91	0.52
1:D:135:ARG:HG3	1:D:136:PHE:HD1	1.75	0.52
1:E:215:PHE:CE1	1:E:237:ARG:HB2	2.45	0.52
1:A:92:ARG:HD3	1:B:78:ASP:O	2.11	0.51
1:C:97:LEU:HB3	1:D:107:LEU:HG	1.91	0.51
1:C:161:LYS:HE3	1:C:174:VAL:CB	2.40	0.51
1:D:53:GLU:HG2	1:D:53:GLU:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:120:MET:O	1:E:124:MET:HG3	2.10	0.51
1:A:167:VAL:HG11	1:A:210:LEU:HD23	1.92	0.51
1:A:163:LEU:HD21	1:A:215:PHE:HZ	1.76	0.51
1:E:167:VAL:HG11	1:E:210:LEU:CD2	2.39	0.51
1:E:33:HIS:HB2	1:E:74:HIS:HB3	1.93	0.51
1:B:202:ASN:ND2	1:B:205:GLN:H	2.08	0.51
1:B:23:ALA:HB2	1:C:65:PRO:HB2	1.93	0.50
1:C:205:GLN:HE21	1:D:189:HIS:CE1	2.30	0.50
1:C:160:ASP:O	1:C:164:ILE:HG12	2.12	0.50
1:C:224:GLY:O	1:C:229:GLN:HG2	2.12	0.50
1:B:85:ARG:HD2	4:C:302:SIN:O3	2.11	0.50
1:B:195:LEU:C	1:B:195:LEU:HD12	2.36	0.50
1:D:68:LEU:C	1:D:68:LEU:HD23	2.37	0.50
1:B:97:LEU:HB3	1:C:107:LEU:HD13	1.94	0.50
1:D:146:ARG:CZ	1:D:192:ARG:HH21	2.24	0.50
1:C:85:ARG:NH2	1:D:82:GLN:HB2	2.26	0.49
1:D:224:GLY:HA3	1:D:229:GLN:HA	1.94	0.49
1:A:191:ILE:HG22	1:A:192:ARG:HG3	1.95	0.49
1:C:215:PHE:CD1	1:C:237:ARG:HB2	2.48	0.49
1:D:207:LEU:O	1:D:211:GLN:HG3	2.13	0.49
1:C:85:ARG:HH22	1:D:83:MET:HE3	1.78	0.48
1:E:28:SER:O	1:E:44:SER:HB2	2.13	0.48
1:B:83:MET:CG	1:B:107:LEU:HG	2.43	0.48
1:C:199:CYS:SG	1:C:201:LEU:HG	2.52	0.48
1:D:209:ARG:O	1:D:213[B]:ARG:HG2	2.13	0.48
1:D:145:VAL:HG13	1:D:233:TYR:HB2	1.96	0.48
1:A:236:ARG:HD3	1:E:211:GLN:NE2	2.28	0.48
1:A:101:ASP:OD1	1:B:103:LYS:HD3	2.13	0.47
1:E:153:GLU:O	1:E:153:GLU:HG3	2.14	0.47
1:A:149:PRO:HG2	1:A:229:GLN:OE1	2.15	0.47
1:A:211:GLN:NE2	1:B:142:CYS:SG	2.88	0.47
1:D:98:ILE:C	1:E:107:LEU:HD11	2.40	0.47
1:E:160:ASP:OD2	1:E:162:SER:HB2	2.14	0.47
1:C:19:ILE:HG23	1:C:20:PRO:HD2	1.96	0.47
1:E:50:LYS:HD2	1:E:91:LEU:O	2.13	0.47
1:C:190:VAL:CG2	1:C:193:PHE:HB2	2.44	0.47
1:B:148:ALA:HA	1:B:149:PRO:HD3	1.77	0.47
1:B:203:SER:O	1:B:206:VAL:HB	2.14	0.47
1:D:188:THR:HG23	1:D:189:HIS:CE1	2.50	0.47
1:D:238:GLU:C	1:D:239:LEU:HD23	2.40	0.46
1:D:153:GLU:OE2	1:D:202:ASN:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:SER:O	1:C:58:ARG:HG3	2.15	0.46
1:D:41:TYR:OH	1:D:81:GLY:O	2.29	0.46
1:D:107:LEU:HD12	1:D:107:LEU:H	1.80	0.46
1:E:213:ARG:O	1:E:237:ARG:NH1	2.48	0.46
1:E:154:ARG:HA	1:E:196:ASN:ND2	2.31	0.46
1:A:156:THR:HB	1:A:192:ARG:HA	1.97	0.46
1:E:127:TRP:CZ2	1:E:131:ARG:HD2	2.51	0.46
1:E:215:PHE:HB3	1:E:235:LEU:HB3	1.97	0.46
1:A:236:ARG:HH11	1:E:211:GLN:NE2	2.14	0.46
1:D:156:THR:HB	1:D:192:ARG:HA	1.98	0.46
1:D:210:LEU:HB3	1:D:235:LEU:HD22	1.97	0.46
1:A:80:ASP:OD2	1:A:83:MET:HG2	2.16	0.46
1:B:212:GLN:O	1:C:137:SER:HA	2.16	0.46
1:D:85[A]:ARG:HH21	4:E:301:SIN:H32	1.80	0.46
1:E:149:PRO:HG2	1:E:229:GLN:HB2	1.97	0.45
1:E:156:THR:HB	1:E:192:ARG:HA	1.97	0.45
1:A:82:GLN:OE1	1:E:85:ARG:NH1	2.49	0.45
1:B:217:ILE:HD12	1:C:218:VAL:HG23	1.98	0.45
1:A:87:ILE:HD13	1:A:120:MET:SD	2.57	0.45
1:B:225:VAL:HG22	1:B:228:SER:CB	2.45	0.45
1:B:235:LEU:HD23	1:B:235:LEU:N	2.31	0.45
1:E:157:LEU:CD1	1:E:195:LEU:HB2	2.47	0.45
1:A:91:LEU:HD21	1:A:117:LEU:HD11	1.99	0.45
1:B:56:ILE:HD11	1:B:60:PHE:HE2	1.81	0.45
1:E:202:ASN:O	1:E:206:VAL:HG23	2.16	0.45
1:A:145:VAL:HG22	1:A:157:LEU:HD23	1.95	0.45
1:A:226:ASP:O	1:A:227:SER:HB2	2.17	0.45
1:A:101:ASP:CG	1:B:103:LYS:HD3	2.42	0.45
1:B:141:GLU:CG	1:B:239:LEU:HD21	2.46	0.45
1:E:195:LEU:C	1:E:195:LEU:HD23	2.42	0.45
1:E:56:ILE:HG23	1:E:57:GLY:N	2.31	0.45
1:D:225:VAL:O	1:D:228:SER:O	2.35	0.44
1:C:47:THR:HG21	1:C:92:ARG:HG2	1.98	0.44
1:C:215:PHE:HB3	1:C:235:LEU:HB3	1.99	0.44
1:C:199:CYS:SG	1:C:200:HIS:N	2.90	0.44
1:D:85[A]:ARG:HB2	1:D:85[A]:ARG:CZ	2.46	0.44
1:A:204:VAL:HG11	1:B:159:GLY:HA2	1.98	0.44
1:B:83:MET:HE2	1:B:83:MET:CA	2.46	0.44
1:C:155:ILE:H	1:C:196:ASN:HB2	1.82	0.44
1:E:139:PRO:HG2	1:E:239:LEU:O	2.17	0.44
1:B:204:VAL:HG12	1:C:144:VAL:CG2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:VAL:HG23	1:B:228:SER:H	1.83	0.44
1:C:81:GLY:H	4:C:302:SIN:C2	2.31	0.44
1:E:122:LEU:O	1:E:126:ARG:HG3	2.18	0.44
1:D:19:ILE:HG23	1:D:20:PRO:HD2	2.00	0.44
1:D:128:LYS:O	1:D:132:GLU:HG3	2.18	0.44
1:C:213:ARG:HB3	1:C:237:ARG:HH21	1.83	0.43
1:E:163:LEU:HD21	1:E:215:PHE:HZ	1.83	0.43
1:D:157:LEU:HD22	1:D:164:ILE:CG1	2.47	0.43
1:B:48:LEU:HD13	1:B:60:PHE:CD1	2.53	0.43
1:B:204:VAL:HG12	1:C:144:VAL:HG23	2.01	0.43
1:C:18:GLY:HA2	1:D:61:ASP:O	2.19	0.43
1:A:191:ILE:HD11	1:E:202:ASN:HB2	2.01	0.43
1:B:28:SER:HB3	1:B:46:ALA:HB2	2.01	0.43
1:B:210:LEU:HD13	1:B:235:LEU:HD12	1.99	0.43
1:C:81:GLY:H	4:C:302:SIN:H22	1.84	0.43
1:D:85[A]:ARG:HH21	4:E:301:SIN:C3	2.32	0.43
1:E:160:ASP:HB2	1:E:189:HIS:CE1	2.53	0.43
1:A:56:ILE:HG23	1:A:57:GLY:H	1.84	0.42
1:D:90:PHE:O	1:D:94:SER:N	2.37	0.42
1:A:215:PHE:CD1	1:A:237:ARG:HB2	2.54	0.42
1:B:23:ALA:HA	1:C:65:PRO:HB3	2.01	0.42
1:C:145:VAL:HG22	1:C:157:LEU:HD11	1.99	0.42
1:E:109:TYR:CE1	1:E:121:LEU:HD22	2.54	0.42
1:D:226:ASP:CG	1:D:227:SER:H	2.28	0.42
1:B:56:ILE:CD1	1:B:60:PHE:HE2	2.32	0.42
1:B:50:LYS:HD2	1:B:91:LEU:O	2.20	0.42
1:D:102:PHE:CZ	1:D:104:ASP:HB2	2.54	0.42
1:D:23:ALA:HB2	1:E:65:PRO:HB2	2.01	0.42
1:A:28:SER:HA	1:A:45:LEU:HB2	2.02	0.41
1:A:78:ASP:C	1:A:79:ARG:HG2	2.45	0.41
1:B:201:LEU:HD12	1:B:205:GLN:NE2	2.34	0.41
1:E:54:SER:OG	1:E:56:ILE:HG22	2.19	0.41
1:C:208:GLU:O	1:C:212:GLN:HG3	2.19	0.41
1:B:105:TYR:CZ	1:B:128:LYS:HE3	2.55	0.41
1:C:152:GLY:O	1:C:153:GLU:HB3	2.20	0.41
1:D:226:ASP:CG	1:D:227:SER:N	2.78	0.41
1:E:58:ARG:HB3	1:E:64:ASP:CG	2.44	0.41
1:A:202:ASN:HB2	1:B:191:ILE:HD11	2.02	0.41
1:B:167:VAL:HG11	1:B:210:LEU:HD23	2.03	0.41
1:C:99:PRO:HD2	1:C:102:PHE:HD1	1.81	0.41
1:B:141:GLU:CD	1:B:239:LEU:HD21	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:146:ARG:NH1	1:D:192:ARG:HH21	2.19	0.41
1:A:47:THR:OG1	1:B:78:ASP:OD2	2.36	0.41
1:B:105:TYR:OH	1:B:128:LYS:HE3	2.21	0.41
1:B:202:ASN:HB2	1:C:191:ILE:HD11	2.03	0.41
1:B:90:PHE:O	1:B:94:SER:N	2.46	0.41
1:B:217:ILE:HG12	1:B:235:LEU:HD22	1.98	0.41
1:D:135:ARG:HG3	1:D:136:PHE:CD1	2.56	0.41
1:E:40:MET:HG3	5:E:302:EDO:C1	2.51	0.41
1:C:42:THR:O	1:D:37:GLY:HA2	2.21	0.41
1:D:138:ARG:NH1	1:D:238:GLU:OE2	2.53	0.41
1:A:134:GLY:O	1:A:138:ARG:HG2	2.21	0.40
1:A:149:PRO:CG	1:B:192:ARG:HH22	2.27	0.40
1:B:113:LYS:HA	1:B:121:LEU:HD11	2.02	0.40
1:B:196:ASN:OD1	1:B:196:ASN:N	2.52	0.40
1:B:199:CYS:SG	1:B:201:LEU:HB2	2.62	0.40
1:C:34:ILE:HG22	1:C:36:VAL:HG23	2.03	0.40
1:C:113:LYS:N	1:C:121:LEU:HD21	2.36	0.40
1:E:102:PHE:CZ	1:E:104:ASP:HB2	2.56	0.40
1:C:50:LYS:HB2	1:C:50:LYS:HE3	1.83	0.40
1:D:160:ASP:OD1	1:D:161:LYS:N	2.54	0.40
1:C:40:MET:HE3	1:C:40:MET:HB3	1.92	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	210/258 (81%)	196 (93%)	14 (7%)	0	100	100
1	B	204/258 (79%)	192 (94%)	12 (6%)	0	100	100
1	C	208/258 (81%)	197 (95%)	11 (5%)	0	100	100
1	D	205/258 (80%)	191 (93%)	14 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	205/258 (80%)	197 (96%)	8 (4%)	0	100	100
All	All	1032/1290 (80%)	973 (94%)	59 (6%)	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	187/234 (80%)	186 (100%)	1 (0%)	86	95
1	B	181/234 (77%)	181 (100%)	0	100	100
1	C	182/234 (78%)	182 (100%)	0	100	100
1	D	180/234 (77%)	180 (100%)	0	100	100
1	E	182/234 (78%)	180 (99%)	2 (1%)	70	87
All	All	912/1170 (78%)	909 (100%)	3 (0%)	91	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	225	VAL
1	E	74	HIS
1	E	237	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	189	HIS
1	A	205	GLN
1	B	74	HIS
1	B	205	GLN
1	C	73	GLN

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Mol	Chain	Res	Type
1	C	74	HIS
1	C	82	GLN
1	C	189	HIS
1	D	189	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 5 are monoatomic - leaving 7 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	$\# Z > 2$
4	SIN	B	301	-	7,7,7	1.02	0	8,8,8	1.44	1 (12%)
4	SIN	D	302	-	7,7,7	1.10	0	8,8,8	1.60	1 (12%)
4	SIN	E	301	-	7,7,7	1.08	0	8,8,8	1.60	1 (12%)
4	SIN	C	302	-	7,7,7	1.01	0	8,8,8	1.88	4 (50%)
5	EDO	E	302	-	3,3,3	0.51	0	2,2,2	0.31	0
5	EDO	B	302	-	3,3,3	0.49	0	2,2,2	0.27	0
4	SIN	A	304	-	7,7,7	1.07	0	8,8,8	1.57	2 (25%)

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	SIN	B	301	-	-	5/5/5/5	-
4	SIN	D	302	-	-	2/5/5/5	-
4	SIN	E	301	-	-	4/5/5/5	-
4	SIN	C	302	-	-	3/5/5/5	-
5	EDO	E	302	-	-	1/1/1/1	-
5	EDO	B	302	-	-	1/1/1/1	-
4	SIN	A	304	-	-	4/5/5/5	-

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	302	SIN	C2-C3-C4	-2.75	107.68	113.60
4	E	301	SIN	C3-C2-C1	-2.40	108.44	113.60
4	D	302	SIN	O2-C1-C2	2.17	121.00	114.03
4	C	302	SIN	O2-C1-C2	2.16	120.97	114.03
4	B	301	SIN	O4-C4-C3	2.15	120.93	114.03
4	A	304	SIN	O2-C1-C2	2.10	120.79	114.03
4	C	302	SIN	C3-C2-C1	-2.07	109.15	113.60
4	A	304	SIN	C3-C2-C1	-2.05	109.20	113.60
4	C	302	SIN	O2-C1-O1	-2.01	118.29	123.30

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	SIN	C1-C2-C3-C4
4	C	302	SIN	C1-C2-C3-C4
5	E	302	EDO	O1-C1-C2-O2
4	B	301	SIN	C2-C3-C4-O3
4	C	302	SIN	O2-C1-C2-C3
4	B	301	SIN	C2-C3-C4-O4
4	C	302	SIN	O1-C1-C2-C3
4	E	301	SIN	C2-C3-C4-O4
4	B	301	SIN	O1-C1-C2-C3
4	B	301	SIN	O2-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
4	A	304	SIN	C2-C3-C4-O4
4	D	302	SIN	O2-C1-C2-C3
4	E	301	SIN	O2-C1-C2-C3
4	E	301	SIN	C2-C3-C4-O3
4	D	302	SIN	O1-C1-C2-C3
5	B	302	EDO	O1-C1-C2-O2
4	A	304	SIN	C2-C3-C4-O3
4	E	301	SIN	O1-C1-C2-C3
4	A	304	SIN	O2-C1-C2-C3
4	A	304	SIN	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301	SIN	1	0
4	E	301	SIN	2	0
4	C	302	SIN	5	0
5	E	302	EDO	2	0
4	A	304	SIN	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	212/258 (82%)	0.21	12 (5%) 30 29	32, 62, 99, 120	2 (0%)
1	B	208/258 (80%)	0.23	10 (4%) 36 35	35, 64, 105, 114	0
1	C	212/258 (82%)	0.36	11 (5%) 34 32	39, 67, 104, 133	0
1	D	207/258 (80%)	0.23	7 (3%) 48 47	27, 62, 105, 127	2 (0%)
1	E	209/258 (81%)	0.22	12 (5%) 30 29	33, 60, 110, 126	0
All	All	1048/1290 (81%)	0.25	52 (4%) 35 34	27, 64, 105, 133	4 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	240	ARG	4.8
1	C	174	VAL	4.6
1	E	187	SER	4.1
1	E	172	GLY	3.9
1	B	187	SER	3.8
1	E	171	ILE	3.7
1	C	173	ASP	3.6
1	B	144	VAL	3.4
1	B	226	ASP	3.3
1	B	142	CYS	3.3
1	B	149	PRO	3.2
1	E	225	VAL	3.1
1	C	232	GLU	3.1
1	A	174	VAL	3.0
1	E	136	PHE	3.0
1	E	150	ASP	2.9
1	D	171	ILE	2.8
1	E	71	LEU	2.8
1	A	159	GLY	2.7
1	B	172	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	135	ARG	2.7
1	D	239	LEU	2.7
1	D	53	GLU	2.7
1	C	192	ARG	2.7
1	C	163	LEU	2.6
1	A	225	VAL	2.5
1	C	197	GLY	2.5
1	B	158	SER	2.5
1	A	150	ASP	2.5
1	A	196	ASN	2.5
1	D	219	GLY	2.5
1	E	161	LYS	2.4
1	A	169	PRO	2.4
1	C	227	SER	2.4
1	C	234	VAL	2.3
1	A	226	ASP	2.3
1	C	225	VAL	2.3
1	E	151	LEU	2.3
1	A	154	ARG	2.2
1	B	198	TYR	2.2
1	D	136	PHE	2.2
1	C	150	ASP	2.2
1	A	187	SER	2.2
1	B	225	VAL	2.2
1	E	56	ILE	2.1
1	D	195	LEU	2.1
1	E	60	PHE	2.1
1	A	221	CYS	2.1
1	B	136	PHE	2.1
1	A	239	LEU	2.1
1	C	151	LEU	2.0
1	E	196	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SIN	C	302	8/8	0.80	0.19	47,59,68,73	0
5	EDO	E	302	4/4	0.82	0.13	69,83,89,99	0
4	SIN	B	301	8/8	0.83	0.14	50,58,75,75	0
4	SIN	E	301	8/8	0.87	0.11	54,61,74,74	0
5	EDO	B	302	4/4	0.87	0.11	68,81,90,91	0
4	SIN	A	304	8/8	0.87	0.11	51,60,72,72	0
4	SIN	D	302	8/8	0.88	0.15	57,67,70,70	0
3	NA	A	302	1/1	0.91	0.22	71,71,71,71	0
3	NA	D	301	1/1	0.96	0.05	50,50,50,50	0
3	NA	A	303	1/1	0.96	0.22	53,53,53,53	0
2	IOD	A	301	1/1	0.99	0.04	58,58,58,58	0
2	IOD	C	301	1/1	0.99	0.06	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.