



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:19 am BST

PDB ID : 5UL7
Title : Structure and function of the divalent anion/Na⁺ symporter from *Vibrio cholerae* and a humanized variant
Authors : Lu, M.
Deposited on : 2017-01-24
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

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

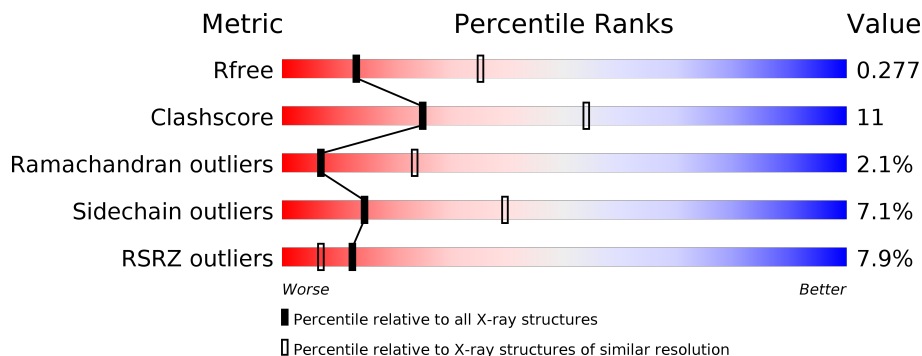
1 Overall quality at a glance

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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 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	445	 8% 74% 22%
1	B	445	 10% 73% 23%
1	C	445	 8% 72% 24%
1	D	445	 6% 71% 25%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13388 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.

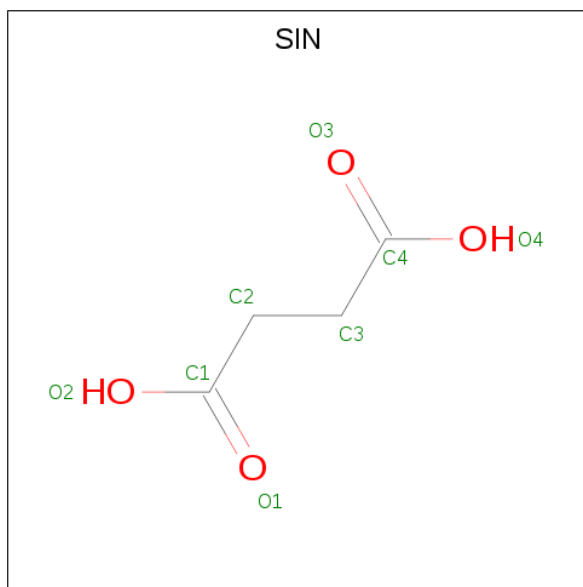
- Molecule 1 is a protein called Transporter, NadC family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	445	3337	2225	522	564	26	0	0	0
1	B	445	3337	2225	522	564	26	0	0	0
1	C	445	3337	2225	522	564	26	0	0	0
1	D	445	3337	2225	522	564	26	0	0	0

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Na	0	0
			2	2		
2	A	2	Total	Na	0	0
			2	2		
2	D	2	Total	Na	0	0
			2	2		
2	C	2	Total	Na	0	0
			2	2		

- Molecule 3 is SUCCINIC ACID (three-letter code: SIN) (formula: C₄H₆O₄).

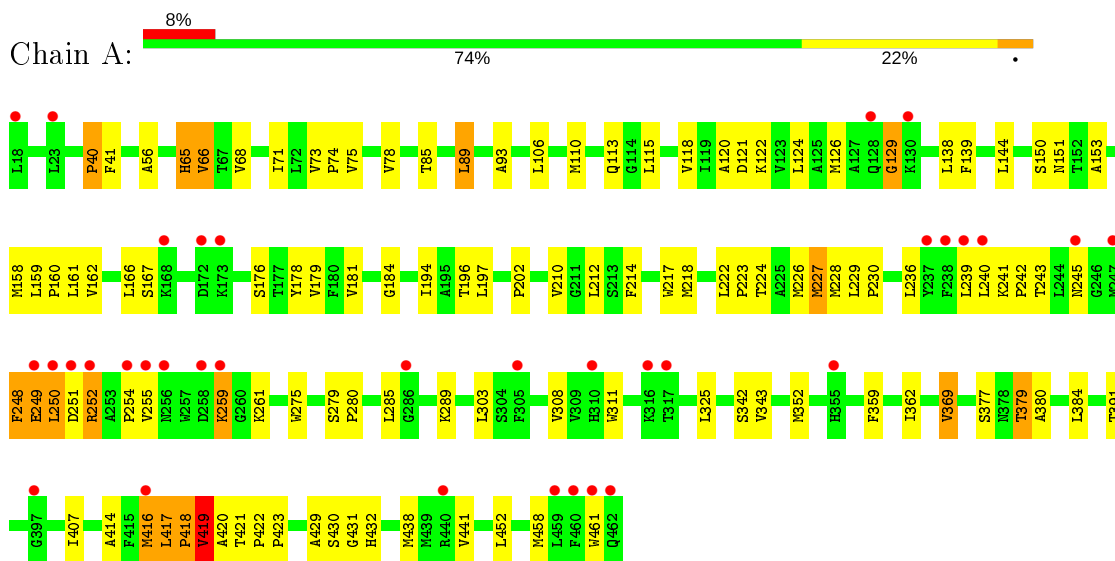


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			8	4 4		
3	B	1	Total	C O	0	0
			8	4 4		
3	C	1	Total	C O	0	0
			8	4 4		
3	D	1	Total	C O	0	0
			8	4 4		

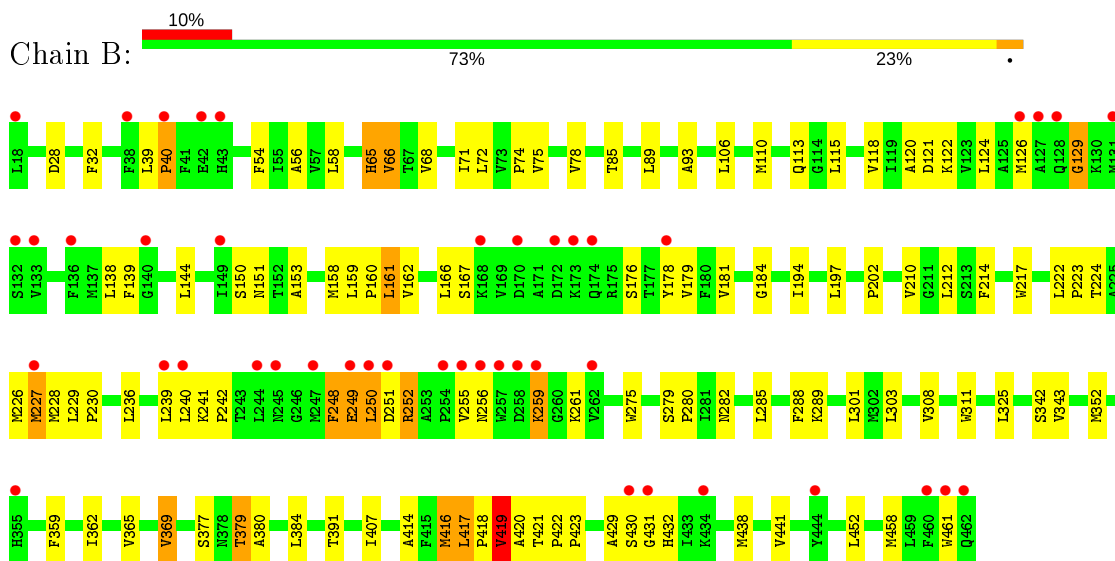
3 Residue-property plots [i](#)

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.

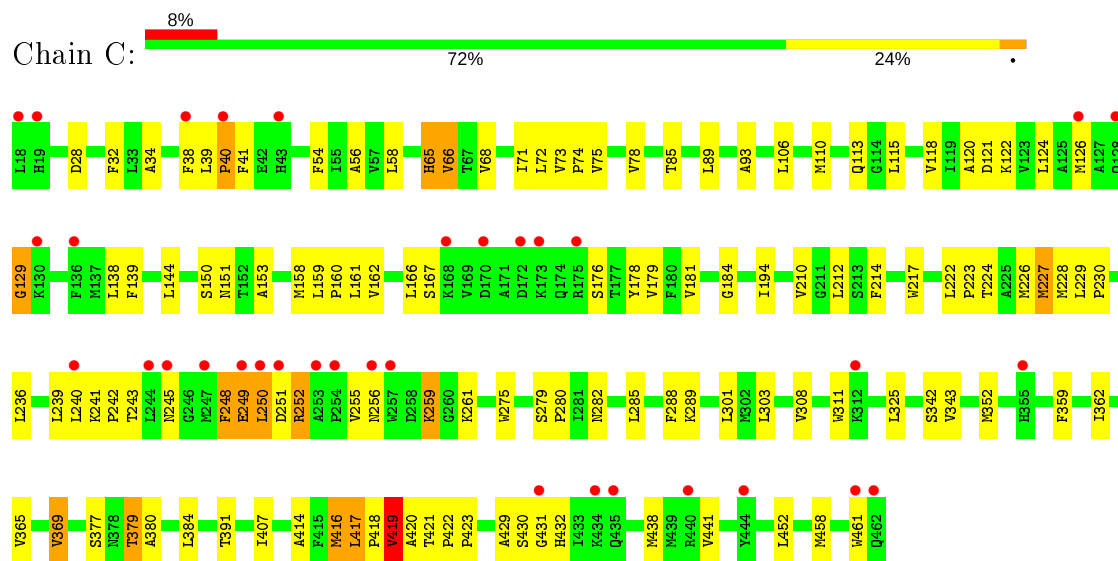
- Molecule 1: Transporter, NadC family



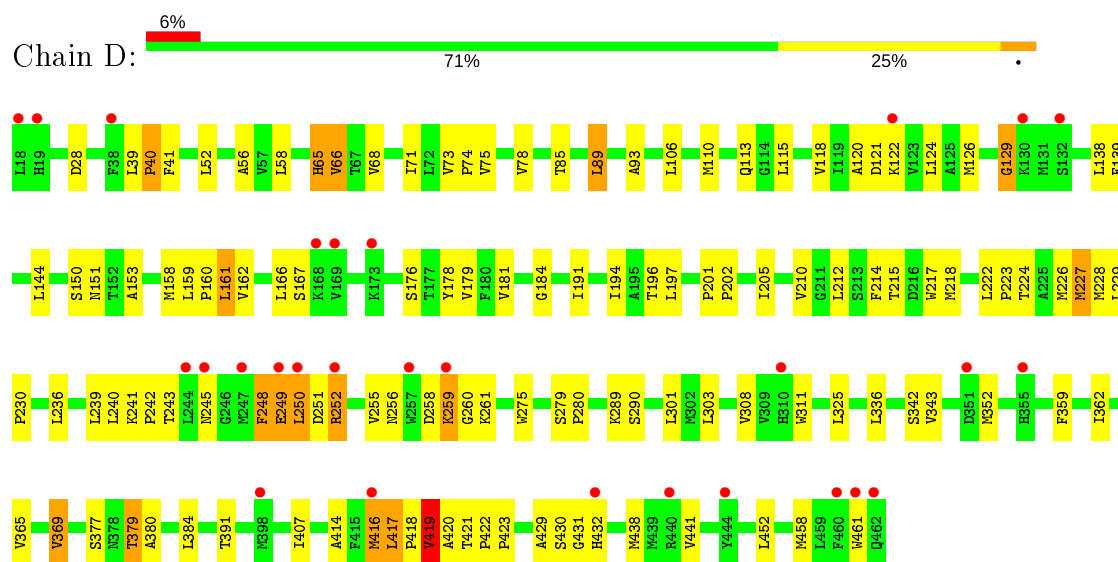
- Molecule 1: Transporter, NadC family



- Molecule 1: Transporter, NadC family



- Molecule 1: Transporter, NadC family



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.69Å 101.92Å 167.75Å 90.00° 98.97° 90.00°	Depositor
Resolution (Å)	15.00 – 2.80 14.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.6 (15.00-2.80) 90.6 (14.98-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.81Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.260 , 0.270 0.269 , 0.277	Depositor DCC
R_{free} test set	3963 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	76.6	Xtrriage
Anisotropy	0.617	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13388	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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: NA, SIN

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	0.35	0/3416	0.51	0/4659
1	B	0.33	0/3416	0.51	0/4659
1	C	0.36	0/3416	0.53	0/4659
1	D	0.37	0/3416	0.53	0/4659
All	All	0.35	0/13664	0.52	0/18636

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	3337	0	3475	74	0
1	B	3337	0	3475	73	0
1	C	3337	0	3475	78	0
1	D	3337	0	3475	86	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	8	0	4	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	4	2	0
3	C	8	0	4	2	0
3	D	8	0	4	2	0
All	All	13388	0	13916	294	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 11.

All (294) 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:D:421:THR:HG22	1:D:423:PRO:HD2	1.14	1.12
1:C:421:THR:HG22	1:C:423:PRO:HD2	1.13	1.11
1:B:421:THR:HG22	1:B:423:PRO:HD2	1.15	1.10
1:A:421:THR:HG22	1:A:423:PRO:HD2	1.10	1.08
1:D:66:VAL:HG12	1:D:325:LEU:HD13	1.61	0.82
1:A:78:VAL:HG11	1:A:85:THR:HG22	1.61	0.82
1:A:421:THR:CG2	1:A:423:PRO:HD2	2.04	0.80
1:B:78:VAL:HG11	1:B:85:THR:HG22	1.61	0.80
1:C:66:VAL:HG12	1:C:325:LEU:HD13	1.62	0.80
1:C:78:VAL:HG11	1:C:85:THR:HG22	1.63	0.79
1:D:78:VAL:HG11	1:D:85:THR:HG22	1.64	0.78
1:A:416:MET:O	1:A:438:MET:HG2	1.85	0.76
1:A:66:VAL:HG12	1:A:325:LEU:HD13	1.68	0.76
1:C:65:HIS:HD2	1:D:311:TRP:HB3	1.49	0.76
1:C:416:MET:O	1:C:438:MET:HG2	1.84	0.76
1:B:416:MET:O	1:B:438:MET:HG2	1.88	0.74
1:D:416:MET:O	1:D:438:MET:HG2	1.88	0.74
1:B:66:VAL:HG12	1:B:325:LEU:HD13	1.67	0.74
1:D:421:THR:CG2	1:D:423:PRO:HD2	2.08	0.74
1:D:362:ILE:HD11	1:D:458:MET:HE2	1.72	0.72
1:B:421:THR:CG2	1:B:423:PRO:HD2	2.09	0.71
1:C:222:LEU:O	1:C:226:MET:HG2	1.90	0.71
1:C:421:THR:CG2	1:C:423:PRO:HD2	2.07	0.71
1:B:222:LEU:O	1:B:226:MET:HG2	1.92	0.70
1:A:122:LYS:O	1:A:126:MET:HG2	1.92	0.69
1:B:122:LYS:O	1:B:126:MET:HG2	1.93	0.68
1:C:362:ILE:HD11	1:C:458:MET:HE2	1.76	0.67
1:B:362:ILE:HD11	1:B:458:MET:HE2	1.76	0.67
1:C:122:LYS:O	1:C:126:MET:HG2	1.95	0.67
1:D:122:LYS:O	1:D:126:MET:HG2	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:VAL:HG11	1:D:301:LEU:HD11	1.78	0.65
1:B:176:SER:HB2	1:B:240:LEU:O	1.96	0.65
1:A:150:SER:HB3	1:A:153:ALA:HB3	1.80	0.63
1:A:311:TRP:HB3	1:B:65:HIS:HD2	1.63	0.63
1:D:222:LEU:O	1:D:226:MET:HG2	1.98	0.63
1:D:176:SER:HB2	1:D:240:LEU:O	1.99	0.63
1:D:418:PRO:C	1:D:420:ALA:H	2.02	0.63
1:A:418:PRO:C	1:A:420:ALA:H	2.02	0.62
1:A:65:HIS:HD2	1:B:311:TRP:HB3	1.64	0.62
1:B:418:PRO:C	1:B:420:ALA:H	2.03	0.62
1:D:150:SER:HB3	1:D:153:ALA:HB3	1.79	0.62
1:A:222:LEU:O	1:A:226:MET:HG2	1.99	0.62
1:A:118:VAL:HG12	1:A:255:VAL:HB	1.82	0.62
1:D:118:VAL:HG12	1:D:255:VAL:HB	1.82	0.62
1:C:414:ALA:HB1	1:C:420:ALA:HB1	1.82	0.61
1:C:65:HIS:HD2	1:D:311:TRP:CB	2.13	0.61
1:A:176:SER:HB2	1:A:240:LEU:O	2.00	0.61
1:C:418:PRO:C	1:C:420:ALA:H	2.04	0.61
1:B:359:PHE:HA	1:B:458:MET:HE1	1.83	0.60
1:C:176:SER:HB2	1:C:240:LEU:O	2.00	0.60
1:C:359:PHE:HA	1:C:458:MET:HE1	1.82	0.60
1:A:421:THR:HG22	1:A:423:PRO:CD	2.06	0.60
1:C:118:VAL:HG12	1:C:255:VAL:HB	1.83	0.59
1:A:414:ALA:HB1	1:A:420:ALA:HB1	1.84	0.59
1:B:414:ALA:HB1	1:B:420:ALA:HB1	1.84	0.59
1:A:359:PHE:CD1	1:A:458:MET:HE3	2.38	0.58
1:B:118:VAL:HG12	1:B:255:VAL:HB	1.84	0.58
1:C:311:TRP:HB3	1:D:65:HIS:HD2	1.68	0.58
1:C:150:SER:HB3	1:C:153:ALA:HB3	1.86	0.57
1:B:159:LEU:HB3	1:B:160:PRO:HD3	1.87	0.57
1:A:167:SER:HB3	1:A:249:GLU:OE2	2.05	0.57
1:C:418:PRO:HD3	1:C:438:MET:CE	2.35	0.57
1:D:359:PHE:HA	1:D:458:MET:HE1	1.87	0.56
1:C:429:ALA:C	1:C:431:GLY:H	2.09	0.56
1:A:429:ALA:C	1:A:431:GLY:H	2.09	0.56
1:D:414:ALA:HB1	1:D:420:ALA:HB1	1.88	0.56
1:D:167:SER:HB3	1:D:249:GLU:OE2	2.06	0.56
1:B:429:ALA:C	1:B:431:GLY:H	2.10	0.55
1:B:150:SER:HB3	1:B:153:ALA:HB3	1.89	0.55
1:C:159:LEU:HB3	1:C:160:PRO:HD3	1.89	0.54
1:A:418:PRO:HD3	1:A:438:MET:CE	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:379:THR:HG23	3:C:503:SIN:O2	2.07	0.54
1:C:167:SER:HB3	1:C:249:GLU:OE2	2.07	0.54
1:D:159:LEU:HB3	1:D:160:PRO:HD3	1.90	0.54
1:B:167:SER:HB3	1:B:249:GLU:OE2	2.08	0.54
1:C:418:PRO:HD3	1:C:438:MET:HE1	1.89	0.53
1:C:71:ILE:O	1:C:74:PRO:HD2	2.09	0.53
1:D:418:PRO:HD3	1:D:438:MET:CE	2.38	0.53
1:A:418:PRO:HD3	1:A:438:MET:HE1	1.90	0.53
1:A:129:GLY:HA2	1:A:248:PHE:HB2	1.91	0.53
1:B:418:PRO:HD3	1:B:438:MET:CE	2.38	0.53
1:A:194:ILE:HG22	1:A:194:ILE:O	2.09	0.53
1:A:311:TRP:CB	1:B:65:HIS:HD2	2.21	0.53
1:D:429:ALA:C	1:D:431:GLY:H	2.10	0.53
1:A:159:LEU:HB3	1:A:160:PRO:HD3	1.91	0.53
1:D:194:ILE:O	1:D:194:ILE:HG22	2.08	0.53
1:B:176:SER:HA	1:B:179:VAL:HG12	1.90	0.53
1:A:71:ILE:O	1:A:74:PRO:HD2	2.09	0.52
1:A:254:PRO:CG	1:C:429:ALA:HB2	2.38	0.52
1:C:212:LEU:HD13	1:C:217:TRP:HD1	1.74	0.52
1:D:138:LEU:HD13	1:D:162:VAL:HG22	1.92	0.52
1:B:212:LEU:HD13	1:B:217:TRP:HD1	1.74	0.52
1:D:377:SER:HB3	1:D:380:ALA:HB3	1.91	0.52
1:A:214:PHE:CD2	1:A:275:TRP:HB3	2.45	0.52
1:C:311:TRP:CB	1:D:65:HIS:HD2	2.23	0.52
1:B:113:GLN:HG3	1:B:261:LYS:HG2	1.92	0.51
1:B:129:GLY:HA2	1:B:248:PHE:HB2	1.92	0.51
1:C:129:GLY:HA2	1:C:248:PHE:HB2	1.91	0.51
1:D:139:PHE:CE2	1:D:184:GLY:HA3	2.46	0.51
1:D:106:LEU:HD12	1:D:303:LEU:HD11	1.92	0.51
1:B:194:ILE:O	1:B:194:ILE:HG22	2.09	0.51
1:C:236:LEU:HD23	1:C:441:VAL:HG11	1.93	0.51
1:A:212:LEU:HD13	1:A:217:TRP:HD1	1.74	0.51
1:D:129:GLY:HA2	1:D:248:PHE:HB2	1.92	0.51
1:A:56:ALA:HA	1:A:384:LEU:HD13	1.92	0.51
1:C:65:HIS:HB3	1:C:68:VAL:HG23	1.92	0.51
1:D:212:LEU:HD13	1:D:217:TRP:HD1	1.76	0.51
1:A:138:LEU:HD13	1:A:162:VAL:HG22	1.93	0.51
1:C:194:ILE:O	1:C:194:ILE:HG22	2.10	0.51
1:C:28:ASP:OD1	1:C:58:LEU:HD13	2.10	0.51
1:A:236:LEU:HD23	1:A:441:VAL:HG11	1.92	0.50
1:A:65:HIS:HD2	1:B:311:TRP:CB	2.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:SER:HA	1:C:179:VAL:HG12	1.93	0.50
1:A:40:PRO:HD2	1:A:41:PHE:CE2	2.47	0.50
1:D:379:THR:HG23	3:D:503:SIN:O1	2.12	0.50
1:A:362:ILE:HD11	1:A:458:MET:HE2	1.92	0.50
1:D:418:PRO:HD3	1:D:438:MET:HE1	1.93	0.50
1:A:176:SER:HA	1:A:179:VAL:HG12	1.94	0.50
1:D:71:ILE:O	1:D:74:PRO:HD2	2.11	0.49
1:D:214:PHE:CD2	1:D:275:TRP:HB3	2.47	0.49
1:D:224:THR:O	1:D:228:MET:HB2	2.12	0.49
1:D:176:SER:HA	1:D:179:VAL:HG12	1.94	0.49
1:B:71:ILE:O	1:B:74:PRO:HD2	2.12	0.49
1:B:369:VAL:HG11	1:B:407:ILE:HG23	1.94	0.49
1:A:139:PHE:CE2	1:A:184:GLY:HA3	2.48	0.49
1:B:418:PRO:O	1:B:419:VAL:HG12	2.13	0.49
1:C:65:HIS:CD2	1:D:311:TRP:HB3	2.39	0.49
1:A:369:VAL:HG11	1:A:407:ILE:HG23	1.94	0.49
1:B:110:MET:HB2	1:B:115:LEU:HB3	1.95	0.49
1:A:377:SER:HB3	1:A:380:ALA:HB3	1.95	0.48
1:C:418:PRO:O	1:C:419:VAL:HG12	2.13	0.48
1:C:422:PRO:HB2	1:C:423:PRO:HD3	1.94	0.48
1:D:113:GLN:HG3	1:D:261:LYS:HG2	1.95	0.48
1:D:197:LEU:HD12	1:D:214:PHE:HA	1.95	0.48
1:A:379:THR:HG23	3:A:503:SIN:O2	2.13	0.48
1:B:222:LEU:HB3	1:B:223:PRO:HD3	1.95	0.48
1:B:422:PRO:HB2	1:B:423:PRO:HD3	1.96	0.48
1:A:113:GLN:HG3	1:A:261:LYS:HG2	1.95	0.48
1:B:139:PHE:CE2	1:B:184:GLY:HA3	2.48	0.48
1:D:236:LEU:HD23	1:D:441:VAL:HG11	1.96	0.48
1:B:377:SER:HB3	1:B:380:ALA:HB3	1.95	0.48
1:D:120:ALA:O	1:D:124:LEU:HB2	2.14	0.48
1:D:56:ALA:HA	1:D:384:LEU:HD13	1.95	0.48
1:D:369:VAL:HG11	1:D:407:ILE:HG23	1.96	0.48
1:B:236:LEU:HD23	1:B:441:VAL:HG11	1.96	0.47
1:C:139:PHE:CE2	1:C:184:GLY:HA3	2.48	0.47
1:D:418:PRO:O	1:D:419:VAL:HG12	2.14	0.47
1:B:379:THR:HG23	3:B:503:SIN:O2	2.14	0.47
1:C:377:SER:HB3	1:C:380:ALA:HB3	1.96	0.47
1:C:214:PHE:CD2	1:C:275:TRP:HB3	2.50	0.47
1:C:113:GLN:HG3	1:C:261:LYS:HG2	1.95	0.47
1:A:75:VAL:HG11	1:B:301:LEU:HD11	1.97	0.47
1:D:40:PRO:HD2	1:D:41:PHE:CE2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:HIS:HB3	1:A:68:VAL:HG23	1.96	0.47
1:C:301:LEU:HD11	1:D:75:VAL:HG11	1.96	0.47
1:A:120:ALA:O	1:A:124:LEU:HB2	2.15	0.47
1:D:65:HIS:HB3	1:D:68:VAL:HG23	1.97	0.46
1:D:118:VAL:HA	1:D:255:VAL:HG21	1.98	0.46
1:A:418:PRO:O	1:A:420:ALA:N	2.44	0.46
1:B:65:HIS:HB3	1:B:68:VAL:HG23	1.97	0.46
1:A:118:VAL:HA	1:A:255:VAL:HG21	1.97	0.46
1:A:422:PRO:HB2	1:A:423:PRO:HD3	1.97	0.46
1:B:214:PHE:CD2	1:B:275:TRP:HB3	2.51	0.46
1:C:166:LEU:HD11	1:C:181:VAL:CB	2.46	0.46
1:D:222:LEU:HB3	1:D:223:PRO:HD3	1.97	0.46
1:C:74:PRO:O	1:C:78:VAL:HG23	2.17	0.45
1:A:418:PRO:C	1:A:420:ALA:N	2.69	0.45
1:C:118:VAL:HA	1:C:255:VAL:HG21	1.97	0.45
1:A:222:LEU:HB3	1:A:223:PRO:HD3	1.99	0.45
1:A:229:LEU:HB3	1:A:230:PRO:HD3	1.99	0.45
1:B:197:LEU:HD12	1:B:214:PHE:HA	1.97	0.45
1:A:417:LEU:HB2	1:A:418:PRO:HD2	1.99	0.45
1:A:418:PRO:O	1:A:419:VAL:HG12	2.16	0.45
1:D:422:PRO:HB2	1:D:423:PRO:HD3	1.97	0.45
1:A:197:LEU:HD12	1:A:214:PHE:HA	1.99	0.45
1:C:138:LEU:HD13	1:C:162:VAL:HG22	1.98	0.45
1:A:65:HIS:ND1	1:A:66:VAL:N	2.65	0.45
1:B:166:LEU:HD11	1:B:181:VAL:CB	2.47	0.45
1:B:138:LEU:HD13	1:B:162:VAL:HG22	1.99	0.44
1:C:342:SER:OG	1:C:391:THR:HB	2.17	0.44
1:A:129:GLY:CA	1:A:248:PHE:HB2	2.47	0.44
1:B:56:ALA:HA	1:B:384:LEU:HD13	1.99	0.44
1:D:342:SER:OG	1:D:391:THR:HB	2.17	0.44
1:D:151:ASN:HB2	3:D:503:SIN:C4	2.48	0.44
1:B:120:ALA:O	1:B:124:LEU:HB2	2.18	0.44
1:B:202:PRO:HD3	1:B:379:THR:HG22	2.00	0.44
1:B:250:LEU:HD23	1:B:250:LEU:H	1.83	0.44
1:D:227:MET:HG2	1:D:452:LEU:HD11	1.99	0.44
1:D:89:LEU:HA	1:D:89:LEU:HD12	1.75	0.44
1:A:224:THR:O	1:A:228:MET:HB2	2.17	0.44
1:B:118:VAL:HA	1:B:255:VAL:HG21	1.99	0.44
1:B:229:LEU:HB3	1:B:230:PRO:HD3	2.00	0.44
1:B:279:SER:HB3	1:B:280:PRO:HD3	1.99	0.44
1:B:417:LEU:HB2	1:B:418:PRO:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:PHE:HB2	1:C:54:PHE:HD1	1.82	0.44
1:C:93:ALA:HB2	1:D:93:ALA:HB2	1.99	0.44
1:C:129:GLY:CA	1:C:248:PHE:HB2	2.48	0.44
1:C:418:PRO:O	1:C:420:ALA:N	2.48	0.44
1:D:229:LEU:HB3	1:D:230:PRO:HD3	1.99	0.44
1:D:418:PRO:C	1:D:420:ALA:N	2.70	0.44
1:A:210:VAL:HG23	1:A:212:LEU:HG	2.00	0.43
1:D:250:LEU:HD23	1:D:250:LEU:H	1.83	0.43
1:A:196:THR:HG22	1:A:218:MET:HG3	2.00	0.43
1:A:250:LEU:HD23	1:A:250:LEU:H	1.83	0.43
1:A:251:ASP:O	1:A:252:ARG:HB3	2.18	0.43
1:B:210:VAL:HG23	1:B:212:LEU:HG	2.00	0.43
1:C:251:ASP:O	1:C:252:ARG:HB3	2.19	0.43
1:A:151:ASN:HB2	3:A:503:SIN:C4	2.49	0.43
1:B:418:PRO:O	1:B:420:ALA:N	2.50	0.43
1:A:202:PRO:HD3	1:A:379:THR:HG22	1.99	0.43
1:B:72:LEU:O	1:B:75:VAL:HG12	2.19	0.43
1:C:210:VAL:HG23	1:C:212:LEU:HG	1.99	0.43
1:A:121:ASP:OD2	1:A:252:ARG:NH2	2.51	0.43
1:C:229:LEU:HB3	1:C:230:PRO:HD3	1.99	0.43
1:D:365:VAL:O	1:D:369:VAL:HG22	2.17	0.43
1:B:418:PRO:HD3	1:B:438:MET:HE3	2.01	0.43
1:C:222:LEU:HB3	1:C:223:PRO:HD3	1.99	0.43
1:D:73:VAL:HB	1:D:74:PRO:HD3	2.01	0.43
1:A:279:SER:HB3	1:A:280:PRO:HD3	2.01	0.43
1:B:32:PHE:HB2	1:B:54:PHE:HD1	1.82	0.43
1:C:121:ASP:OD2	1:C:252:ARG:NH2	2.52	0.43
1:C:279:SER:HB3	1:C:280:PRO:HD3	1.99	0.43
1:C:369:VAL:HG11	1:C:407:ILE:HG23	2.01	0.43
1:D:166:LEU:HD11	1:D:181:VAL:CB	2.49	0.43
1:A:106:LEU:HD12	1:A:303:LEU:HD11	2.00	0.43
1:B:418:PRO:HD3	1:B:438:MET:HE1	1.99	0.43
1:C:110:MET:HB2	1:C:115:LEU:HB3	2.01	0.43
1:D:52:LEU:HD13	1:D:336:LEU:HD21	2.00	0.43
1:A:89:LEU:HD12	1:A:89:LEU:HA	1.83	0.42
1:B:342:SER:OG	1:B:391:THR:HB	2.19	0.42
1:C:56:ALA:HA	1:C:384:LEU:HD13	2.01	0.42
1:D:202:PRO:HD3	1:D:379:THR:HG22	2.00	0.42
1:D:417:LEU:HB2	1:D:418:PRO:HD2	2.01	0.42
1:B:106:LEU:HD12	1:B:303:LEU:HD11	2.00	0.42
1:C:418:PRO:C	1:C:420:ALA:N	2.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:417:LEU:HB2	1:C:418:PRO:HD2	2.00	0.42
1:D:418:PRO:O	1:D:420:ALA:N	2.48	0.42
1:A:166:LEU:HD11	1:A:181:VAL:CB	2.48	0.42
1:C:250:LEU:HD23	1:C:250:LEU:H	1.84	0.42
1:D:196:THR:HG22	1:D:218:MET:HG3	2.01	0.42
1:A:110:MET:HB2	1:A:115:LEU:HB3	2.00	0.42
1:B:227:MET:HG2	1:B:452:LEU:HD11	2.01	0.42
1:D:251:ASP:O	1:D:252:ARG:HB3	2.20	0.42
1:A:93:ALA:HB2	1:B:93:ALA:HB2	2.02	0.42
1:C:34:ALA:O	1:C:38:PHE:HB2	2.20	0.42
1:C:72:LEU:O	1:C:75:VAL:HG12	2.19	0.42
1:C:75:VAL:HG11	1:D:301:LEU:CD1	2.49	0.42
1:D:279:SER:HB3	1:D:280:PRO:HD3	2.01	0.42
1:B:251:ASP:O	1:B:252:ARG:HB3	2.19	0.42
1:B:121:ASP:OD2	1:B:252:ARG:NH2	2.52	0.42
1:B:151:ASN:HB2	3:B:503:SIN:C4	2.49	0.42
1:C:120:ALA:O	1:C:124:LEU:HB2	2.18	0.42
1:A:359:PHE:HA	1:A:458:MET:HE1	2.00	0.42
1:A:227:MET:HG2	1:A:452:LEU:HD11	2.00	0.42
1:C:151:ASN:HB2	3:C:503:SIN:C4	2.50	0.42
1:D:201:PRO:O	1:D:205:ILE:HG13	2.19	0.42
1:B:28:ASP:OD1	1:B:58:LEU:HD13	2.19	0.42
1:C:243:THR:C	1:C:245:ASN:H	2.23	0.41
1:D:129:GLY:CA	1:D:248:PHE:HB2	2.49	0.41
1:D:258:ASP:O	1:D:260:GLY:N	2.53	0.41
1:B:161:LEU:HD11	1:B:252:ARG:NH2	2.35	0.41
1:C:282:ASN:OD1	1:C:288:PHE:N	2.53	0.41
1:C:106:LEU:HD12	1:C:303:LEU:HD11	2.01	0.41
1:C:227:MET:HG2	1:C:452:LEU:HD11	2.02	0.41
1:C:365:VAL:O	1:C:369:VAL:HG22	2.20	0.41
1:D:39:LEU:HA	1:D:40:PRO:HD3	1.87	0.41
1:D:248:PHE:HD2	1:D:249:GLU:H	1.69	0.41
1:D:121:ASP:OD2	1:D:252:ARG:NH2	2.52	0.41
1:D:65:HIS:ND1	1:D:66:VAL:N	2.69	0.41
1:A:342:SER:OG	1:A:391:THR:HB	2.20	0.41
1:B:224:THR:O	1:B:228:MET:HB2	2.21	0.41
1:D:194:ILE:O	1:D:194:ILE:CG2	2.69	0.41
1:B:282:ASN:OD1	1:B:288:PHE:N	2.54	0.41
1:C:85:THR:HG21	1:D:290:SER:H	1.85	0.41
1:D:227:MET:HB3	1:D:452:LEU:HD21	2.02	0.41
1:B:365:VAL:O	1:B:369:VAL:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:ILE:CD1	1:B:458:MET:HE2	2.48	0.41
1:A:73:VAL:HB	1:A:74:PRO:HD3	2.03	0.41
1:C:224:THR:O	1:C:228:MET:HB2	2.20	0.41
1:C:73:VAL:HB	1:C:74:PRO:HD3	2.03	0.41
1:D:161:LEU:HD11	1:D:252:ARG:HH22	1.86	0.41
1:A:243:THR:C	1:A:245:ASN:H	2.24	0.40
1:D:191:ILE:HG12	1:D:228:MET:CE	2.51	0.40
1:D:210:VAL:HG23	1:D:212:LEU:HG	2.02	0.40
1:D:243:THR:C	1:D:245:ASN:H	2.24	0.40
1:D:28:ASP:OD1	1:D:58:LEU:HD13	2.21	0.40
1:B:248:PHE:HD2	1:B:249:GLU:H	1.69	0.40
1:D:110:MET:HB2	1:D:115:LEU:HB3	2.02	0.40
1:B:65:HIS:ND1	1:B:66:VAL:N	2.69	0.40
1:C:39:LEU:HA	1:C:40:PRO:HD3	1.86	0.40
1:A:227:MET:HB3	1:A:452:LEU:HD21	2.04	0.40
1:B:39:LEU:HA	1:B:40:PRO:HD3	1.86	0.40
1:C:40:PRO:HD2	1:C:41:PHE:CE2	2.57	0.40
1:D:161:LEU:HD11	1:D:252:ARG:NH2	2.36	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	443/445 (100%)	404 (91%)	29 (6%)	10 (2%)	6	21
1	B	443/445 (100%)	402 (91%)	32 (7%)	9 (2%)	7	24
1	C	443/445 (100%)	404 (91%)	30 (7%)	9 (2%)	7	24
1	D	443/445 (100%)	403 (91%)	31 (7%)	9 (2%)	7	24
All	All	1772/1780 (100%)	1613 (91%)	122 (7%)	37 (2%)	7	23

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	PRO
1	A	419	VAL
1	B	242	PRO
1	B	419	VAL
1	C	242	PRO
1	C	419	VAL
1	D	242	PRO
1	D	259	LYS
1	D	419	VAL
1	A	252	ARG
1	A	259	LYS
1	A	430	SER
1	B	252	ARG
1	B	259	LYS
1	B	430	SER
1	C	66	VAL
1	C	252	ARG
1	C	259	LYS
1	C	430	SER
1	D	40	PRO
1	D	252	ARG
1	A	40	PRO
1	A	129	GLY
1	A	416	MET
1	B	40	PRO
1	B	129	GLY
1	B	416	MET
1	C	40	PRO
1	C	129	GLY
1	C	416	MET
1	D	129	GLY
1	D	416	MET
1	D	430	SER
1	B	66	VAL
1	D	66	VAL
1	A	66	VAL
1	A	418	PRO

5.3.2 Protein sidechains [i](#)

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	351/353 (99%)	327 (93%)	24 (7%)	16	42
1	B	351/353 (99%)	326 (93%)	25 (7%)	14	39
1	C	351/353 (99%)	326 (93%)	25 (7%)	14	39
1	D	351/353 (99%)	326 (93%)	25 (7%)	14	39
All	All	1404/1412 (99%)	1305 (93%)	99 (7%)	14	39

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	89	LEU
1	A	144	LEU
1	A	158	MET
1	A	161	LEU
1	A	178	TYR
1	A	227	MET
1	A	239	LEU
1	A	241	LYS
1	A	248	PHE
1	A	249	GLU
1	A	250	LEU
1	A	259	LYS
1	A	285	LEU
1	A	289	LYS
1	A	308	VAL
1	A	343	VAL
1	A	352	MET
1	A	369	VAL
1	A	379	THR
1	A	417	LEU
1	A	419	VAL
1	A	432	HIS
1	A	461	TRP
1	B	65	HIS
1	B	89	LEU
1	B	144	LEU
1	B	158	MET
1	B	161	LEU
1	B	178	TYR

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Mol	Chain	Res	Type
1	B	227	MET
1	B	239	LEU
1	B	241	LYS
1	B	248	PHE
1	B	249	GLU
1	B	250	LEU
1	B	256	ASN
1	B	259	LYS
1	B	285	LEU
1	B	289	LYS
1	B	308	VAL
1	B	343	VAL
1	B	352	MET
1	B	369	VAL
1	B	379	THR
1	B	417	LEU
1	B	419	VAL
1	B	432	HIS
1	B	461	TRP
1	C	65	HIS
1	C	89	LEU
1	C	144	LEU
1	C	158	MET
1	C	161	LEU
1	C	178	TYR
1	C	227	MET
1	C	239	LEU
1	C	241	LYS
1	C	248	PHE
1	C	249	GLU
1	C	250	LEU
1	C	256	ASN
1	C	259	LYS
1	C	285	LEU
1	C	289	LYS
1	C	308	VAL
1	C	343	VAL
1	C	352	MET
1	C	369	VAL
1	C	379	THR
1	C	417	LEU
1	C	419	VAL

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Mol	Chain	Res	Type
1	C	432	HIS
1	C	461	TRP
1	D	65	HIS
1	D	89	LEU
1	D	144	LEU
1	D	158	MET
1	D	161	LEU
1	D	178	TYR
1	D	215	THR
1	D	227	MET
1	D	239	LEU
1	D	241	LYS
1	D	248	PHE
1	D	249	GLU
1	D	250	LEU
1	D	256	ASN
1	D	259	LYS
1	D	289	LYS
1	D	308	VAL
1	D	343	VAL
1	D	352	MET
1	D	369	VAL
1	D	379	THR
1	D	417	LEU
1	D	419	VAL
1	D	432	HIS
1	D	461	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	432	HIS
1	B	43	HIS
1	B	44	ASN
1	B	151	ASN
1	B	432	HIS
1	C	43	HIS
1	C	65	HIS
1	C	151	ASN
1	C	432	HIS
1	D	43	HIS

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Mol	Chain	Res	Type
1	D	432	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	$\# Z > 2$
3	SIN	B	503	-	1,7,7	0.11	0	2,8,8	1.61	0
3	SIN	A	503	-	1,7,7	0.08	0	2,8,8	1.63	0
3	SIN	D	503	-	1,7,7	0.06	0	2,8,8	1.61	0
3	SIN	C	503	-	1,7,7	0.11	0	2,8,8	1.88	1 (50%)

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
3	SIN	B	503	-	-	0/1/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SIN	A	503	-	-	0/1/5/5	-
3	SIN	D	503	-	-	0/1/5/5	-
3	SIN	C	503	-	-	0/1/5/5	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	503	SIN	C3-C2-C1	-2.34	108.74	112.67

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	503	SIN	2	0
3	A	503	SIN	2	0
3	D	503	SIN	2	0
3	C	503	SIN	2	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	445/445 (100%)	0.07	35 (7%) 12 7	42, 84, 135, 189	0
1	B	445/445 (100%)	0.23	44 (9%) 7 4	57, 97, 157, 196	0
1	C	445/445 (100%)	0.04	34 (7%) 13 7	42, 77, 140, 195	0
1	D	445/445 (100%)	-0.00	28 (6%) 20 12	25, 74, 124, 176	0
All	All	1780/1780 (100%)	0.09	141 (7%) 12 7	25, 83, 146, 196	0

All (141) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	462	GLN	11.4
1	B	462	GLN	8.7
1	B	245	ASN	8.1
1	B	133	VAL	7.7
1	C	250	LEU	7.7
1	C	462	GLN	7.5
1	C	355	HIS	6.9
1	A	18	LEU	6.7
1	B	168	LYS	6.3
1	A	128	GLN	6.2
1	C	254	PRO	6.2
1	A	130	LYS	5.8
1	B	355	HIS	5.8
1	C	18	LEU	5.5
1	B	132	SER	5.4
1	A	168	LYS	5.3
1	A	460	PHE	5.2
1	D	168	LYS	5.2
1	D	130	LYS	5.1
1	D	245	ASN	5.0
1	C	245	ASN	5.0

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Mol	Chain	Res	Type	RSRZ
1	B	240	LEU	4.9
1	B	126	MET	4.7
1	B	18	LEU	4.6
1	B	257	TRP	4.6
1	A	462	GLN	4.6
1	A	245	ASN	4.5
1	B	255	VAL	4.5
1	A	172	ASP	4.3
1	C	128	GLN	4.3
1	A	240	LEU	4.3
1	A	249	GLU	4.2
1	A	256	ASN	4.2
1	A	255	VAL	4.2
1	D	460	PHE	4.1
1	A	355	HIS	4.1
1	A	286	GLY	4.1
1	C	38	PHE	4.0
1	D	259	LYS	4.0
1	C	253	ALA	4.0
1	C	130	LYS	3.9
1	B	461	TRP	3.9
1	D	247	MET	3.9
1	B	259	LYS	3.9
1	D	440	ARG	3.8
1	C	168	LYS	3.8
1	D	250	LEU	3.8
1	C	126	MET	3.8
1	D	432	HIS	3.8
1	B	256	ASN	3.7
1	D	444	TYR	3.7
1	B	247	MET	3.7
1	C	257	TRP	3.6
1	C	249	GLU	3.6
1	B	40	PRO	3.6
1	C	19	HIS	3.6
1	B	128	GLN	3.5
1	A	397	GLY	3.5
1	A	461	TRP	3.4
1	C	256	ASN	3.4
1	B	434	LYS	3.4
1	D	173	LYS	3.4
1	A	254	PRO	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	40	PRO	3.3
1	C	172	ASP	3.3
1	B	258	ASP	3.3
1	B	430	SER	3.2
1	C	461	TRP	3.2
1	C	247	MET	3.2
1	A	310	HIS	3.1
1	B	173	LYS	3.1
1	D	18	LEU	3.1
1	B	254	PRO	3.1
1	C	251	ASP	3.1
1	A	258	ASP	3.1
1	B	136	PHE	3.0
1	B	172	ASP	3.0
1	B	178	TYR	3.0
1	B	250	LEU	3.0
1	A	252	ARG	3.0
1	B	170	ASP	3.0
1	D	249	GLU	2.9
1	C	431	GLY	2.9
1	C	444	TYR	2.9
1	C	240	LEU	2.8
1	D	169	VAL	2.8
1	D	132	SER	2.8
1	D	355	HIS	2.8
1	A	317	THR	2.7
1	D	252	ARG	2.7
1	B	244	LEU	2.7
1	B	262	VAL	2.7
1	D	351	ASP	2.6
1	B	460	PHE	2.6
1	A	251	ASP	2.6
1	A	316	LYS	2.6
1	B	140	GLY	2.6
1	B	239	LEU	2.6
1	B	127	ALA	2.5
1	A	416	MET	2.5
1	B	227	MET	2.5
1	D	416	MET	2.5
1	A	250	LEU	2.5
1	C	43	HIS	2.4
1	C	173	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	305	PHE	2.4
1	A	23	LEU	2.4
1	A	459	LEU	2.4
1	C	170	ASP	2.4
1	D	398	MET	2.4
1	B	38	PHE	2.4
1	D	38	PHE	2.4
1	C	434	LYS	2.3
1	C	244	LEU	2.3
1	A	238	PHE	2.3
1	B	174	GLN	2.3
1	A	247	MET	2.3
1	B	444	TYR	2.3
1	C	440	ARG	2.3
1	A	259	LYS	2.3
1	D	310	HIS	2.3
1	B	43	HIS	2.2
1	B	431	GLY	2.2
1	B	42	GLU	2.2
1	C	136	PHE	2.2
1	D	244	LEU	2.2
1	D	257	TRP	2.2
1	B	251	ASP	2.1
1	C	175	ARG	2.1
1	D	122	LYS	2.1
1	A	440	ARG	2.1
1	B	249	GLU	2.1
1	B	149	ILE	2.1
1	A	239	LEU	2.1
1	D	461	TRP	2.1
1	C	435	GLN	2.0
1	C	312	LYS	2.0
1	D	19	HIS	2.0
1	A	173	LYS	2.0
1	A	237	TYR	2.0
1	B	131	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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. 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
3	SIN	B	503	8/8	0.72	0.40	90,94,98,99	0
2	NA	B	501	1/1	0.80	0.26	99,99,99,99	0
2	NA	D	502	1/1	0.80	0.28	79,79,79,79	0
2	NA	B	502	1/1	0.81	0.20	87,87,87,87	0
2	NA	A	502	1/1	0.81	0.26	64,64,64,64	0
3	SIN	D	503	8/8	0.85	0.23	93,96,99,100	0
2	NA	D	501	1/1	0.89	0.41	79,79,79,79	0
3	SIN	A	503	8/8	0.90	0.32	76,78,80,80	0
3	SIN	C	503	8/8	0.91	0.38	78,81,83,85	0
2	NA	C	501	1/1	0.91	0.09	89,89,89,89	0
2	NA	C	502	1/1	0.94	0.21	74,74,74,74	0
2	NA	A	501	1/1	0.96	0.10	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.