



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:52 AM GMT

PDB ID : 3CHG
Title : The compatible solute-binding protein OpuAC from *Bacillus subtilis* in complex with DMSA
Authors : Smits, S.H.J.; Hoing, M.; Lecher, J.; Jebbar, M.; Schmitt, L.; Bremer, E.
Deposited on : 2008-03-09
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

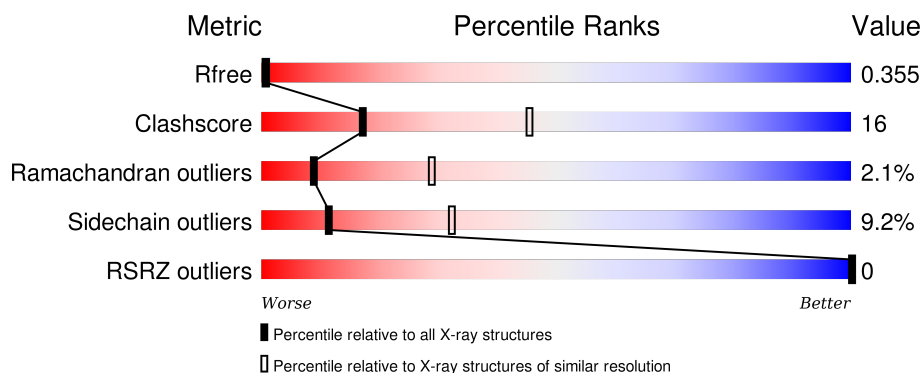
1 Overall quality at a glance

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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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

Mol	Chain	Length	Quality of chain
1	A	268	 67% 28% . .
1	B	268	 71% 22% . .
1	C	268	 62% 28% 7% .
1	D	268	 56% 33% 6% . .

2 Entry composition [i](#)

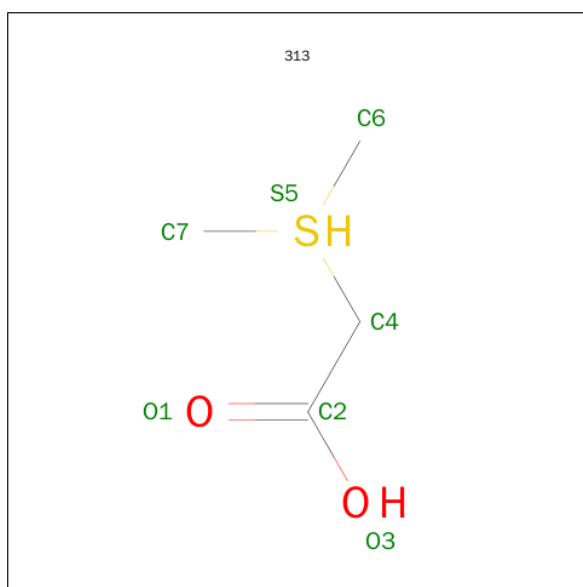
There are 2 unique types of molecules in this entry. The entry contains 8104 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycine betaine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	257	Total	C	N	O	S	0	1	0
			2019	1290	331	387	11			
1	A	258	Total	C	N	O	S	0	0	0
			2019	1289	332	388	10			
1	B	258	Total	C	N	O	S	0	0	0
			2019	1289	332	388	10			
1	C	258	Total	C	N	O	S	0	0	0
			2019	1289	332	388	10			

- Molecule 2 is (DIMETHYL-LAMBDA 4 -SULFANYL)ACETIC ACID (three-letter code: 313) (formula: C₄H₁₀O₂S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	O	S	0	0
			7	4	2	1		
2	B	1	Total	C	O	S	0	0
			7	4	2	1		

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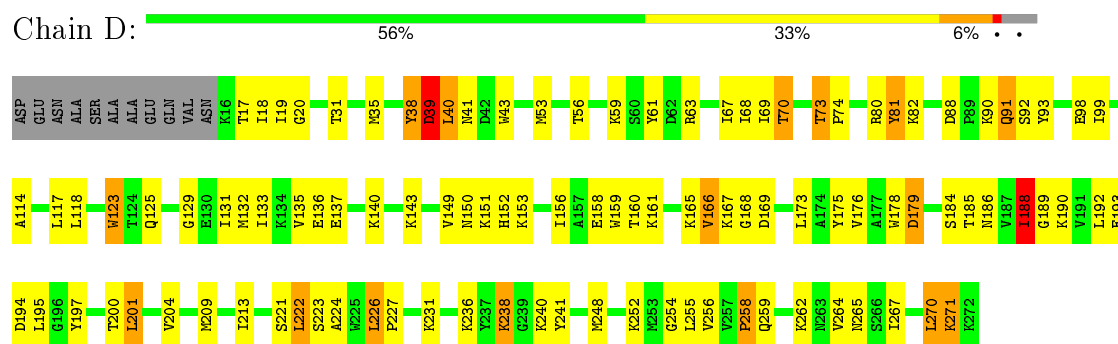
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	O	S	0	0
			7	4	2	1		
2	A	1	Total	C	O	S	0	0
			7	4	2	1		

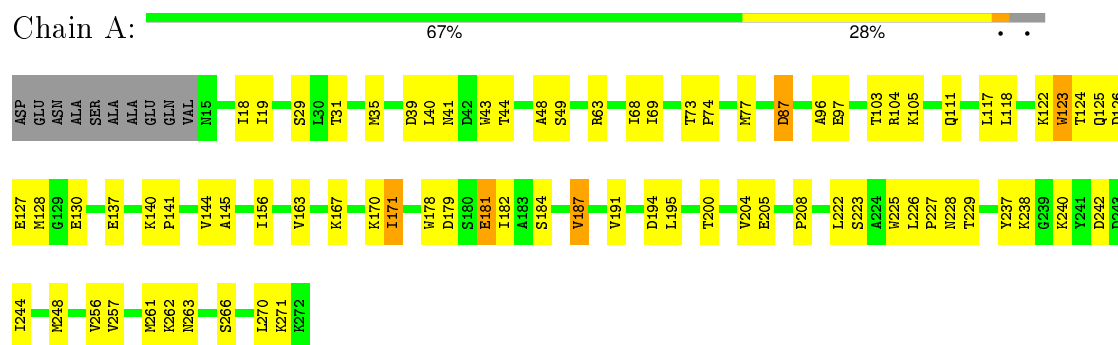
3 Residue-property plots

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

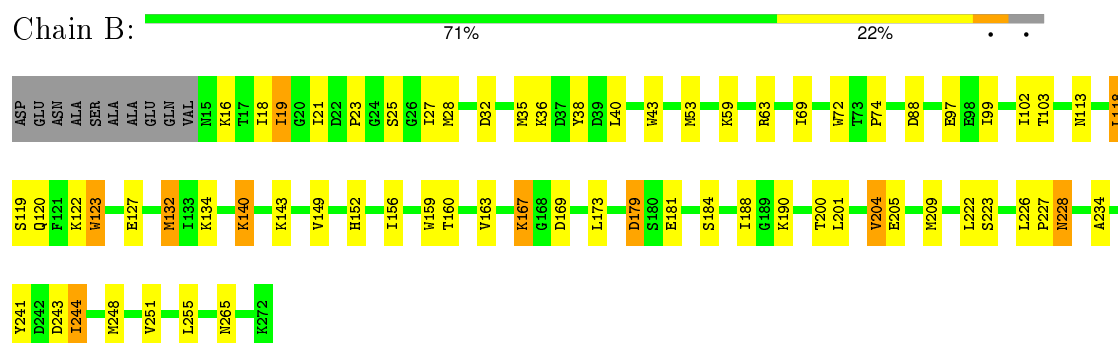
- Molecule 1: Glycine betaine-binding protein



- Molecule 1: Glycine betaine-binding protein

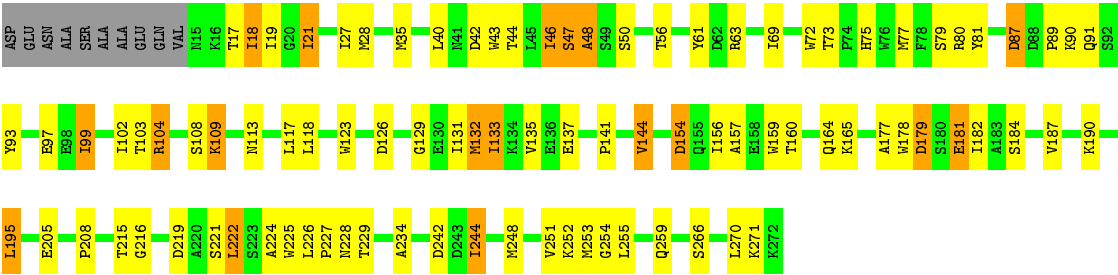


- Molecule 1: Glycine betaine-binding protein



- Molecule 1: Glycine betaine-binding protein

Chain C: 62% 28% 7% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.51Å 150.61Å 58.96Å 90.00° 104.54° 90.00°	Depositor
Resolution (Å)	19.56 – 2.80 19.51 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.0 (19.56-2.80) 93.0 (19.51-2.80)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.285 , 0.359 0.289 , 0.355	Depositor DCC
R_{free} test set	1119 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 21.1	EDS
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 21785 reflections	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	8104	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 80.10 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.8216e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
313

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	0.39	0/2062	0.55	0/2786
1	B	0.37	0/2062	0.53	0/2786
1	C	0.36	0/2062	0.53	1/2786 (0.0%)
1	D	0.58	0/2062	0.55	1/2785 (0.0%)
All	All	0.43	0/8248	0.54	2/11143 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	3
1	D	0	10
All	All	0	13

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	226	LEU	C-N-CD	-6.08	107.23	120.60
1	C	195	LEU	CA-CB-CG	5.19	127.24	115.30

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	132	MET	Peptide
1	C	46	ILE	Peptide

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Mol	Chain	Res	Type	Group
1	C	47	SER	Peptide
1	D	168	GLY	Peptide
1	D	188	ILE	Peptide
1	D	200	THR	Peptide
1	D	201	LEU	Peptide
1	D	258	PRO	Peptide
1	D	270	LEU	Peptide
1	D	271	LYS	Peptide
1	D	38	TYR	Peptide
1	D	39	ASP	Peptide
1	D	40	LEU	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2019	0	2034	37	1
1	B	2019	0	2034	39	0
1	C	2019	0	2034	63	0
1	D	2019	0	2036	126	2
2	A	7	0	9	0	0
2	B	7	0	9	1	0
2	C	7	0	9	0	0
2	D	7	0	9	2	0
All	All	8104	0	8174	263	2

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 16.

All (263) 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:114:ALA:O	1:D:118:LEU:HD13	1.19	1.33
1:C:132:MET:O	1:C:133:ILE:HG13	1.36	1.21
1:D:270:LEU:HD13	1:D:270:LEU:O	1.44	1.17
1:D:74:PRO:HB3	1:D:125:GLN:HG3	1.17	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:114:ALA:O	1:D:118:LEU:CD1	1.94	1.15
1:D:150:ASN:O	1:D:153:LYS:HD2	1.56	1.05
1:D:188:ILE:HD13	1:D:222:LEU:HD21	1.38	1.03
1:D:166:VAL:HG11	1:D:197:TYR:CE1	1.96	1.00
1:D:74:PRO:HB3	1:D:125:GLN:CG	1.93	0.99
1:C:132:MET:O	1:C:133:ILE:CG1	2.14	0.95
1:D:18:ILE:HD13	1:D:68:ILE:HB	1.49	0.94
1:D:166:VAL:HG11	1:D:197:TYR:CD1	2.05	0.91
1:D:40:LEU:HD23	1:D:271:LYS:HD2	1.52	0.90
1:D:150:ASN:O	1:D:153:LYS:CD	2.19	0.89
1:D:184:SER:O	1:D:188:ILE:CG2	2.23	0.86
1:D:166:VAL:CG1	1:D:197:TYR:CE1	2.57	0.86
1:C:21:ILE:O	1:C:47:SER:O	1.92	0.86
1:D:184:SER:O	1:D:188:ILE:HG22	1.75	0.86
1:D:20:GLY:HA2	1:D:70:THR:OG1	1.75	0.85
1:D:74:PRO:CB	1:D:125:GLN:HG3	2.05	0.84
1:D:188:ILE:HD13	1:D:222:LEU:CD2	2.08	0.83
1:D:166:VAL:CG1	1:D:197:TYR:CD1	2.61	0.82
1:A:184:SER:O	1:A:187:VAL:HG12	1.81	0.81
1:B:99:ILE:CG2	1:B:248:MET:HB2	2.11	0.81
1:A:237:TYR:HD1	1:A:240:LYS:HD2	1.45	0.81
1:C:28:MET:HE1	1:C:46:ILE:O	1.81	0.80
1:D:18:ILE:HD13	1:D:68:ILE:CB	2.12	0.79
1:D:59:LYS:HA	1:D:63:ARG:HH21	1.48	0.79
1:D:118:LEU:CD1	1:D:118:LEU:H	1.97	0.77
1:C:73:THR:CG2	1:C:252:LYS:HB2	2.16	0.76
1:C:18:ILE:HD11	1:C:255:LEU:CD2	2.16	0.75
1:C:46:ILE:HG22	1:C:47:SER:H	1.50	0.75
1:C:102:ILE:HD11	1:C:221:SER:HB3	1.69	0.74
1:D:213:ILE:HD11	1:D:221:SER:HB2	1.68	0.74
1:C:18:ILE:HD11	1:C:255:LEU:HD22	1.68	0.74
1:D:61:TYR:CE1	1:D:81:TYR:HB3	2.22	0.74
1:D:38:TYR:O	1:D:39:ASP:HB2	1.85	0.74
1:C:73:THR:HG22	1:C:252:LYS:HB2	1.70	0.73
1:D:270:LEU:CD1	1:D:270:LEU:O	2.30	0.73
1:D:129:GLY:O	1:D:133:ILE:HG13	1.88	0.73
1:D:188:ILE:CD1	1:D:222:LEU:HD21	2.16	0.73
1:D:222:LEU:N	1:D:222:LEU:HD12	2.02	0.73
1:D:226:LEU:HB3	1:D:227:PRO:HA	1.70	0.73
1:B:99:ILE:HG22	1:B:248:MET:HB2	1.70	0.72
1:D:118:LEU:N	1:D:118:LEU:HD12	2.04	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:40:LEU:O	1:D:43:TRP:HB2	1.90	0.71
1:C:17:THR:HG22	1:C:18:ILE:H	1.56	0.71
1:C:108:SER:O	1:C:109:LYS:HB2	1.90	0.70
1:D:166:VAL:CG2	1:D:167:LYS:N	2.54	0.70
1:D:158:GLU:HG2	1:D:161:LYS:NZ	2.07	0.70
1:C:113:ASN:HD21	1:C:164:GLN:H	1.39	0.70
1:D:43:TRP:CZ2	1:D:270:LEU:HD13	2.27	0.69
1:D:149:VAL:O	1:D:153:LYS:HG2	1.92	0.69
1:D:18:ILE:HD12	1:D:255:LEU:HD22	1.74	0.69
1:D:193:GLU:O	1:D:195:LEU:N	2.26	0.68
1:A:87:ASP:HB2	1:A:266:SER:HA	1.73	0.68
1:D:173:LEU:HB2	1:D:201:LEU:HD12	1.76	0.68
1:D:43:TRP:CZ2	1:D:270:LEU:CD1	2.77	0.68
1:A:237:TYR:CD1	1:A:240:LYS:HD2	2.28	0.68
1:D:59:LYS:HA	1:D:63:ARG:NH2	2.07	0.68
1:D:173:LEU:HB2	1:D:201:LEU:CD1	2.23	0.68
1:D:188:ILE:O	1:D:192:LEU:HB2	1.94	0.67
1:D:114:ALA:C	1:D:118:LEU:HD13	2.10	0.67
1:D:150:ASN:C	1:D:153:LYS:HD2	2.14	0.67
1:D:73:THR:HG22	1:D:252:LYS:H	1.60	0.67
1:A:48:ALA:O	1:A:49:SER:HB3	1.94	0.67
1:D:186:ASN:ND2	1:D:201:LEU:HD21	2.10	0.66
1:C:141:PRO:HA	1:C:144:VAL:HG12	1.77	0.66
1:D:118:LEU:CD1	1:D:118:LEU:N	2.58	0.66
1:D:140:LYS:HB2	1:D:143:LYS:HD2	1.78	0.66
1:D:18:ILE:HD13	1:D:68:ILE:CG2	2.27	0.65
1:D:99:ILE:HA	1:D:224:ALA:O	1.95	0.65
1:D:61:TYR:CD2	1:D:81:TYR:HD1	2.15	0.65
1:D:222:LEU:CD1	1:D:222:LEU:N	2.61	0.63
1:B:59:LYS:O	1:B:59:LYS:HG2	1.97	0.63
1:C:117:LEU:HD11	1:C:160:THR:HG22	1.80	0.63
1:D:18:ILE:CD1	1:D:68:ILE:CG2	2.77	0.63
1:D:136:GLU:O	1:D:136:GLU:HG3	1.99	0.62
1:A:171:ILE:O	1:A:171:ILE:HG23	1.99	0.62
1:B:181:GLU:HG2	1:B:223:SER:HA	1.82	0.62
1:D:166:VAL:HG22	1:D:167:LYS:N	2.15	0.61
1:D:88:ASP:OD2	1:D:91:GLN:HA	2.00	0.61
1:A:68:ILE:HG13	1:A:257:VAL:HG12	1.81	0.61
1:D:209:MET:SD	1:D:221:SER:OG	2.58	0.60
1:D:61:TYR:CZ	1:D:81:TYR:HB3	2.36	0.60
1:D:18:ILE:CD1	1:D:68:ILE:HG21	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:ASP:CG	1:D:91:GLN:HA	2.21	0.60
1:D:193:GLU:C	1:D:195:LEU:H	2.04	0.59
1:C:132:MET:O	1:C:133:ILE:CB	2.49	0.59
1:B:205:GLU:O	1:B:209:MET:HB2	2.03	0.59
1:C:99:ILE:CG2	1:C:248:MET:HB3	2.32	0.59
1:C:102:ILE:CD1	1:C:221:SER:HB3	2.33	0.59
1:C:21:ILE:HG22	1:C:50:SER:HA	1.85	0.59
1:D:117:LEU:HG	1:D:160:THR:HG22	1.85	0.59
1:B:102:ILE:HG22	1:B:243:ASP:HA	1.84	0.58
1:D:132[B]:MET:HE3	1:D:179:ASP:HA	1.83	0.58
1:D:166:VAL:HG11	1:D:197:TYR:CZ	2.36	0.58
1:C:35:MET:HA	1:C:40:LEU:HB2	1.85	0.58
1:C:117:LEU:HD13	1:C:195:LEU:HD11	1.85	0.57
1:D:131:ILE:O	1:D:135:VAL:HG23	2.05	0.57
1:B:35:MET:HA	1:B:40:LEU:HD12	1.86	0.57
1:D:40:LEU:CD2	1:D:271:LYS:HD2	2.31	0.57
1:C:99:ILE:HD11	1:C:222:LEU:O	2.05	0.57
1:D:150:ASN:HA	1:D:153:LYS:HD2	1.87	0.57
1:A:69:ILE:HG12	1:A:256:VAL:O	2.03	0.57
1:D:166:VAL:HG12	1:D:197:TYR:CE1	2.38	0.56
1:C:97:GLU:HB3	1:C:251:VAL:HG23	1.86	0.56
1:C:154:ASP:N	1:C:154:ASP:OD2	2.36	0.56
1:D:43:TRP:CH2	1:D:270:LEU:HD11	2.40	0.56
1:D:118:LEU:H	1:D:118:LEU:HD12	1.63	0.56
1:D:150:ASN:CA	1:D:153:LYS:HD2	2.36	0.56
1:C:225:TRP:HB3	1:C:229:THR:CG2	2.36	0.55
1:D:158:GLU:HG2	1:D:161:LYS:HZ2	1.72	0.55
1:D:213:ILE:CD1	1:D:221:SER:HB2	2.36	0.55
1:C:87:ASP:O	1:C:89:PRO:HD3	2.07	0.55
1:B:149:VAL:HG23	1:B:156:ILE:HG12	1.89	0.55
1:A:35:MET:HA	1:A:40:LEU:HB2	1.87	0.54
1:A:141:PRO:HA	1:A:144:VAL:HG12	1.87	0.54
1:D:40:LEU:HD11	1:D:267:ILE:HG22	1.90	0.54
1:D:73:THR:HG23	1:D:125:GLN:OE1	2.07	0.54
1:D:43:TRP:CZ2	1:D:270:LEU:HD11	2.42	0.54
1:B:156:ILE:HA	1:B:159:TRP:HB2	1.89	0.54
1:D:188:ILE:O	1:D:192:LEU:HD12	2.07	0.54
1:A:97:GLU:OE2	1:A:229:THR:HG22	2.08	0.54
1:B:226:LEU:HD21	1:B:234:ALA:HB2	1.90	0.54
1:C:18:ILE:HD11	1:C:255:LEU:HD21	1.90	0.53
1:A:156:ILE:H	1:A:156:ILE:HD12	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:TRP:HE1	1:A:271:LYS:HD3	1.74	0.53
1:D:188:ILE:HG21	1:D:222:LEU:CD2	2.39	0.53
1:D:166:VAL:HG23	1:D:167:LYS:H	1.73	0.53
1:A:226:LEU:HB3	1:A:227:PRO:HA	1.90	0.53
1:D:90:LYS:O	1:D:91:GLN:C	2.47	0.52
1:D:150:ASN:C	1:D:153:LYS:HG3	2.30	0.52
1:B:103:THR:HG23	1:B:244:ILE:HD12	1.91	0.52
1:A:43:TRP:HZ2	1:A:270:LEU:O	1.92	0.52
1:A:68:ILE:HG13	1:A:257:VAL:CG1	2.40	0.51
1:D:193:GLU:C	1:D:195:LEU:N	2.62	0.51
1:D:118:LEU:H	1:D:118:LEU:HD13	1.74	0.51
1:A:178:TRP:O	1:A:182:ILE:HG13	2.10	0.51
1:D:149:VAL:HG23	1:D:156:ILE:HG12	1.93	0.51
1:D:166:VAL:CG2	1:D:167:LYS:H	2.23	0.51
1:D:188:ILE:HG21	1:D:222:LEU:HD23	1.92	0.51
1:C:43:TRP:CZ2	1:C:270:LEU:HB3	2.46	0.51
1:A:181:GLU:HG2	1:A:223:SER:HA	1.93	0.50
1:C:103:THR:OG1	1:C:242:ASP:HB2	2.11	0.50
1:C:46:ILE:HG22	1:C:47:SER:N	2.22	0.50
1:C:69:ILE:O	1:C:255:LEU:HA	2.11	0.50
1:C:226:LEU:HD11	1:C:234:ALA:HB2	1.93	0.50
1:A:117:LEU:HD12	1:A:163:VAL:HG11	1.94	0.50
1:B:32:ASP:O	1:B:36:LYS:HD2	2.11	0.49
1:C:47:SER:O	1:C:48:ALA:O	2.30	0.49
1:C:19:ILE:HD13	1:C:56:THR:HG21	1.94	0.49
1:B:113:ASN:ND2	1:B:163:VAL:HG13	2.28	0.49
1:B:113:ASN:HD21	1:B:163:VAL:HG13	1.77	0.49
1:B:27:ILE:HG22	2:B:2:313:O3	2.12	0.49
1:C:184:SER:O	1:C:187:VAL:HG12	2.12	0.49
1:B:127:GLU:HG2	1:B:152:HIS:HE1	1.77	0.49
1:C:93:TYR:CE1	1:C:254:GLY:HA2	2.48	0.49
1:B:241:TYR:HE1	1:B:243:ASP:HB2	1.79	0.48
1:B:16:LYS:HB3	1:B:43:TRP:CD2	2.49	0.48
1:D:150:ASN:HA	1:D:153:LYS:CD	2.44	0.48
1:D:176:VAL:HG23	1:D:209:MET:HG3	1.95	0.48
1:D:61:TYR:CD2	1:D:81:TYR:CD1	3.00	0.48
1:D:59:LYS:HG2	1:D:63:ARG:HH21	1.79	0.48
1:C:132:MET:C	1:C:133:ILE:HG13	2.24	0.47
1:C:117:LEU:CD1	1:C:195:LEU:HD11	2.43	0.47
1:C:72:TRP:HA	1:C:253:MET:HA	1.96	0.47
1:D:82:LYS:O	1:D:259:GLN:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:GLN:HG2	1:A:248:MET:SD	2.54	0.47
1:A:127:GLU:O	1:A:130:GLU:HB2	2.14	0.47
1:C:79:SER:O	1:C:80:ARG:HG2	2.15	0.47
1:B:226:LEU:HB3	1:B:227:PRO:HA	1.97	0.47
1:D:226:LEU:HB3	1:D:227:PRO:CA	2.42	0.47
1:D:175:TYR:HB3	1:D:185:THR:HG21	1.96	0.47
1:B:127:GLU:HG2	1:B:152:HIS:CE1	2.49	0.47
1:D:238:LYS:C	1:D:240:LYS:H	2.17	0.47
1:D:38:TYR:OH	1:D:88:ASP:OD1	2.25	0.47
1:D:133:ILE:O	1:D:137:GLU:HG2	2.15	0.46
1:C:75:HIS:ND1	1:C:77:MET:HB2	2.30	0.46
1:A:31:THR:O	1:A:35:MET:HG3	2.16	0.46
1:A:103:THR:OG1	1:A:242:ASP:HB2	2.14	0.46
1:D:150:ASN:C	1:D:153:LYS:CG	2.84	0.46
1:D:262:LYS:O	1:D:265:ASN:ND2	2.45	0.46
1:A:39:ASP:HA	1:A:41:ASN:HD22	1.81	0.46
1:D:43:TRP:CZ2	1:D:270:LEU:O	2.69	0.46
1:D:39:ASP:O	1:D:271:LYS:CE	2.64	0.46
1:C:226:LEU:HB3	1:C:227:PRO:HA	1.98	0.46
1:D:186:ASN:ND2	1:D:201:LEU:CD2	2.78	0.46
1:B:25:SER:OG	1:B:28:MET:HB2	2.15	0.46
1:D:123:TRP:CZ2	1:D:184:SER:HB3	2.50	0.46
1:C:27:ILE:HD12	1:C:253:MET:HB2	1.97	0.46
1:B:74:PRO:HD3	1:B:251:VAL:HG12	1.97	0.46
1:D:90:LYS:O	1:D:92:SER:N	2.50	0.45
1:D:158:GLU:HG2	1:D:161:LYS:HZ1	1.78	0.45
1:D:80:ARG:HB3	1:D:81:TYR:CE2	2.52	0.45
1:D:123:TRP:HZ2	1:D:184:SER:HB3	1.81	0.45
1:D:123:TRP:O	1:D:248:MET:HG2	2.17	0.45
1:B:97:GLU:OE1	1:B:228:ASN:HB3	2.17	0.45
1:A:178:TRP:O	1:A:182:ILE:N	2.50	0.45
1:C:87:ASP:HB2	1:C:266:SER:HA	1.99	0.45
1:C:61:TYR:CE1	1:C:81:TYR:HB3	2.52	0.44
1:D:93:TYR:CE1	1:D:254:GLY:HA2	2.52	0.44
1:C:156:ILE:HA	1:C:159:TRP:HB2	1.98	0.44
1:D:238:LYS:O	1:D:240:LYS:N	2.46	0.44
1:D:69:ILE:HG13	1:D:256:VAL:HB	2.00	0.44
1:D:166:VAL:HG21	1:D:169:ASP:HB2	1.99	0.44
1:A:205:GLU:HB2	1:A:208:PRO:HG2	2.00	0.44
1:C:87:ASP:C	1:C:89:PRO:HD3	2.39	0.43
1:A:122:LYS:O	1:A:123:TRP:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:SER:O	1:C:109:LYS:CB	2.62	0.43
1:C:225:TRP:HB3	1:C:229:THR:HG21	2.01	0.43
1:B:132:MET:HE3	1:B:179:ASP:HA	1.99	0.43
1:D:59:LYS:CA	1:D:63:ARG:HH21	2.26	0.43
1:C:205:GLU:HB2	1:C:208:PRO:HG2	2.00	0.43
1:D:178:TRP:CD2	2:D:1:313:H4A	2.53	0.43
1:C:129:GLY:O	1:C:133:ILE:HD12	2.19	0.43
1:B:19:ILE:HG23	1:B:69:ILE:HG22	1.99	0.43
1:B:21:ILE:HG21	1:B:72:TRP:HH2	1.84	0.43
1:A:96:ALA:HB2	1:C:48:ALA:N	2.34	0.43
1:B:134:LYS:HE3	1:C:63:ARG:HG2	2.01	0.43
1:C:102:ILE:HG12	1:C:224:ALA:HB2	2.01	0.42
1:C:225:TRP:HB3	1:C:229:THR:HG23	2.01	0.42
1:C:131:ILE:O	1:C:135:VAL:HG23	2.19	0.42
1:A:145:ALA:HB1	1:A:187:VAL:HA	2.01	0.42
1:A:225:TRP:HB3	1:A:229:THR:CG2	2.49	0.42
1:C:90:LYS:O	1:C:91:GLN:HB2	2.20	0.42
1:C:179:ASP:HA	1:C:182:ILE:HD12	2.02	0.42
1:A:191:VAL:O	1:A:195:LEU:HG	2.19	0.42
1:D:150:ASN:C	1:D:153:LYS:CD	2.83	0.42
1:B:59:LYS:HA	1:B:63:ARG:NH2	2.35	0.42
1:A:170:LYS:O	1:A:171:ILE:HB	2.18	0.42
1:A:126:ASP:OD1	1:A:127:GLU:N	2.52	0.42
1:D:158:GLU:C	1:D:160:THR:N	2.73	0.42
1:D:39:ASP:O	1:D:271:LYS:HE3	2.19	0.42
1:D:19:ILE:HG22	1:D:53:MET:HG2	2.02	0.42
1:A:74:PRO:HG2	1:A:128:MET:SD	2.60	0.42
1:B:38:TYR:OH	1:B:88:ASP:OD2	2.29	0.42
1:D:158:GLU:O	1:D:161:LYS:HG2	2.20	0.41
1:C:104:ARG:HB2	1:C:104:ARG:HH11	1.84	0.41
1:A:105:LYS:HD2	1:A:105:LYS:HA	1.88	0.41
1:B:140:LYS:HB2	1:B:143:LYS:HG3	2.01	0.41
1:B:204:VAL:HB	1:B:205:GLU:H	1.59	0.41
1:D:67:ILE:HG23	1:D:258:PRO:HD3	2.02	0.41
1:B:122:LYS:O	1:B:123:TRP:HB3	2.20	0.41
1:D:59:LYS:O	1:D:63:ARG:NE	2.53	0.41
1:C:154:ASP:O	1:C:157:ALA:HB3	2.21	0.41
1:D:186:ASN:HD21	1:D:201:LEU:CD2	2.33	0.41
1:D:178:TRP:CE2	2:D:1:313:H4A	2.56	0.41
1:B:19:ILE:HD11	1:B:53:MET:HA	2.03	0.41
1:B:27:ILE:HG12	1:B:27:ILE:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:TRP:HZ2	1:B:184:SER:HB3	1.86	0.41
1:A:261:MET:C	1:A:263:ASN:H	2.23	0.41
1:B:173:LEU:HB2	1:B:201:LEU:HD23	2.02	0.41
1:A:111:GLN:HE22	1:A:171:ILE:HB	1.86	0.40
1:B:118:LEU:O	1:B:120:GLN:N	2.54	0.40
1:C:178:TRP:HB2	1:C:181:GLU:HB2	2.02	0.40
1:D:192:LEU:O	1:D:193:GLU:C	2.60	0.40
1:D:132[B]:MET:HE3	1:D:179:ASP:CA	2.48	0.40
1:C:226:LEU:CB	1:C:227:PRO:HA	2.50	0.40
1:B:188:ILE:HA	1:B:188:ILE:HD13	2.00	0.40
1:D:31:THR:O	1:D:35:MET:HG3	2.21	0.40
1:B:167:LYS:HA	1:B:167:LYS:HD3	1.78	0.40
1:C:244:ILE:HD13	1:C:244:ILE:H	1.85	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:41:ASN:ND2	1:D:165:LYS:N[1_655]	2.05	0.15
1:D:236:LYS:NZ	1:A:104:ARG:NH2[2_656]	2.10	0.10

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	256/268 (96%)	226 (88%)	27 (10%)	3 (1%)	16	47
1	B	256/268 (96%)	235 (92%)	19 (7%)	2 (1%)	24	58
1	C	256/268 (96%)	219 (86%)	29 (11%)	8 (3%)	5	17
1	D	256/268 (96%)	216 (84%)	32 (12%)	8 (3%)	5	17
All	All	1024/1072 (96%)	896 (88%)	107 (10%)	21 (2%)	9	29

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	39	ASP
1	A	171	ILE
1	C	109	LYS
1	C	133	ILE
1	D	91	GLN
1	D	189	GLY
1	D	194	ASP
1	D	238	LYS
1	C	18	ILE
1	A	262	LYS
1	D	152	HIS
1	C	216	GLY
1	D	231	LYS
1	A	238	LYS
1	B	119	SER
1	C	177	ALA
1	C	215	THR
1	D	159	TRP
1	C	48	ALA
1	B	204	VAL
1	C	21	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	216/223 (97%)	193 (89%)	23 (11%)	8	24
1	B	216/223 (97%)	198 (92%)	18 (8%)	14	38
1	C	216/223 (97%)	195 (90%)	21 (10%)	10	29
1	D	216/223 (97%)	199 (92%)	17 (8%)	15	40
All	All	864/892 (97%)	785 (91%)	79 (9%)	11	33

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	17	THR
1	D	56	THR
1	D	70	THR
1	D	73	THR
1	D	81	TYR
1	D	98	GLU
1	D	123	TRP
1	D	151	LYS
1	D	166	VAL
1	D	179	ASP
1	D	188	ILE
1	D	190	LYS
1	D	204	VAL
1	D	222	LEU
1	D	223	SER
1	D	241	TYR
1	D	264	VAL
1	A	18	ILE
1	A	19	ILE
1	A	29	SER
1	A	44	THR
1	A	63	ARG
1	A	73	THR
1	A	77	MET
1	A	87	ASP
1	A	118	LEU
1	A	123	TRP
1	A	124	THR
1	A	137	GLU
1	A	140	LYS
1	A	167	LYS
1	A	179	ASP
1	A	181	GLU
1	A	187	VAL
1	A	194	ASP
1	A	200	THR
1	A	204	VAL
1	A	222	LEU
1	A	228	ASN
1	A	244	ILE
1	B	18	ILE
1	B	19	ILE
1	B	23	PRO

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Mol	Chain	Res	Type
1	B	118	LEU
1	B	123	TRP
1	B	132	MET
1	B	140	LYS
1	B	160	THR
1	B	167	LYS
1	B	169	ASP
1	B	179	ASP
1	B	190	LYS
1	B	200	THR
1	B	222	LEU
1	B	228	ASN
1	B	244	ILE
1	B	255	LEU
1	B	265	ASN
1	C	42	ASP
1	C	44	THR
1	C	87	ASP
1	C	99	ILE
1	C	104	ARG
1	C	118	LEU
1	C	123	TRP
1	C	126	ASP
1	C	137	GLU
1	C	144	VAL
1	C	154	ASP
1	C	165	LYS
1	C	179	ASP
1	C	181	GLU
1	C	190	LYS
1	C	219	ASP
1	C	222	LEU
1	C	228	ASN
1	C	244	ILE
1	C	259	GLN
1	C	271	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	91	GLN
1	D	100	HIS

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Mol	Chain	Res	Type
1	D	113	ASN
1	D	120	GLN
1	D	152	HIS
1	D	186	ASN
1	D	259	GLN
1	A	15	ASN
1	A	41	ASN
1	A	111	GLN
1	A	113	ASN
1	A	186	ASN
1	A	228	ASN
1	B	113	ASN
1	B	186	ASN
1	B	228	ASN
1	C	113	ASN
1	C	120	GLN
1	C	186	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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
2	313	A	4	-	2,6,6	6.10	2 (100%)	3,7,7	6.11	3 (100%)
2	313	B	2	-	2,6,6	5.96	2 (100%)	3,7,7	6.41	3 (100%)
2	313	C	3	-	2,6,6	5.94	2 (100%)	3,7,7	5.24	3 (100%)
2	313	D	1	-	2,6,6	6.09	2 (100%)	3,7,7	5.41	3 (100%)

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	313	A	4	-	-	0/0/4/4	0/0/0/0
2	313	B	2	-	-	0/0/4/4	0/0/0/0
2	313	C	3	-	-	0/0/4/4	0/0/0/0
2	313	D	1	-	-	0/0/4/4	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	4	313	C6-S5	-6.17	1.66	1.78
2	D	1	313	C7-S5	-6.10	1.66	1.78
2	D	1	313	C6-S5	-6.08	1.66	1.78
2	A	4	313	C7-S5	-6.03	1.66	1.78
2	B	2	313	C7-S5	-6.02	1.66	1.78
2	C	3	313	C7-S5	-6.01	1.66	1.78
2	B	2	313	C6-S5	-5.91	1.66	1.78
2	C	3	313	C6-S5	-5.88	1.66	1.78

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	313	C7-S5-C4	2.77	105.96	101.56
2	D	1	313	C6-S5-C4	3.97	107.87	101.56
2	A	4	313	C6-S5-C4	4.26	108.33	101.56
2	C	3	313	C6-S5-C4	4.94	109.40	101.56
2	D	1	313	C7-S5-C4	5.21	109.83	101.56
2	A	4	313	C7-S5-C4	6.04	111.16	101.56
2	B	2	313	C6-S5-C7	6.09	108.95	101.63
2	B	2	313	C6-S5-C4	6.52	111.92	101.56
2	B	2	313	C7-S5-C4	6.62	112.08	101.56
2	D	1	313	C6-S5-C7	6.71	109.70	101.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	313	C6-S5-C7	7.11	110.17	101.63
2	A	4	313	C6-S5-C7	7.56	110.72	101.63

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	313	1	0
2	D	1	313	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	258/268 (96%)	-0.06	0 100 100	10, 16, 21, 24	0
1	B	258/268 (96%)	-0.00	0 100 100	9, 16, 22, 23	0
1	C	258/268 (96%)	-0.04	0 100 100	9, 15, 21, 23	0
1	D	257/268 (95%)	0.07	0 100 100	8, 16, 20, 22	0
All	All	1031/1072 (96%)	-0.01	0 100 100	8, 16, 21, 24	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	313	D	1	7/7	0.93	0.18	0.12	28,28,28,28	0
2	313	B	2	7/7	0.93	0.20	-0.02	18,18,19,19	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	313	A	4	7/7	0.95	0.18	-0.32	27,28,28,28	0
2	313	C	3	7/7	0.96	0.15	-1.29	16,16,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.