



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:13 AM GMT

PDB ID : 2ZZD
Title : Recombinant thiocyanate hydrolase, air-oxidized form of holo-enzyme
Authors : Arakawa, T.; Kawano, Y.; Katayama, Y.; Yohda, M.; Odaka, M.
Deposited on : 2009-02-09
Resolution : 1.78 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

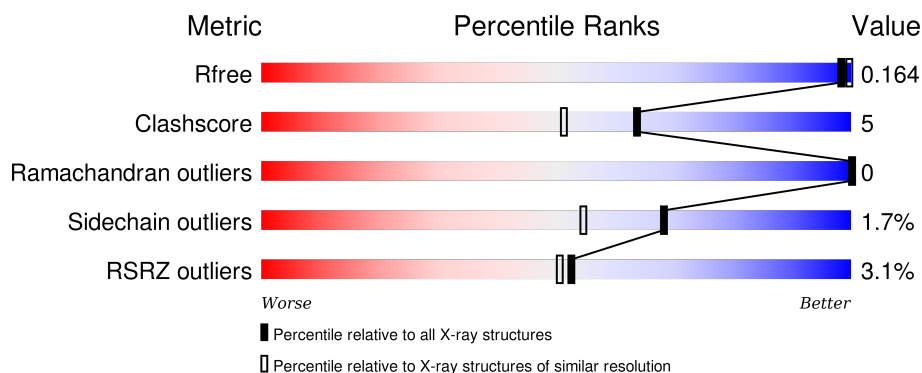
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	6655 (1.80-1.76)
Clashscore	102246	7658 (1.80-1.76)
Ramachandran outliers	100387	7570 (1.80-1.76)
Sidechain outliers	100360	7569 (1.80-1.76)
RSRZ outliers	91569	6671 (1.80-1.76)

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	126	<div> <div>4%</div> <div>85%</div> <div>9%</div> <div>6%</div> </div>
1	D	126	<div> <div>2%</div> <div>87%</div> <div>6%</div> <div>5%</div> </div>
1	G	126	<div> <div>4%</div> <div>86%</div> <div>9%</div> <div>5%</div> </div>
1	J	126	<div> <div>2%</div> <div>89%</div> <div>6%</div> <div>5%</div> </div>
2	B	157	<div> <div>0%</div> <div>85%</div> <div>11%</div> <div>••</div> </div>

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Mol	Chain	Length	Quality of chain
2	E	157	
2	H	157	
2	K	157	
3	C	243	
3	F	243	
3	I	243	
3	L	243	

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	FRU	B	6002	-	-	-	X
4	FRU	L	6001	-	-	-	X
4	FRU	L	6003	-	-	-	X
7	BGC	F	5001	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 19243 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 Thiocyanate hydrolase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	0	0
			965	614	160	187	4			
1	D	120	Total	C	N	O	S	0	0	0
			974	620	162	188	4			
1	G	120	Total	C	N	O	S	0	0	0
			974	620	162	188	4			
1	J	120	Total	C	N	O	S	0	0	0
			974	620	162	188	4			

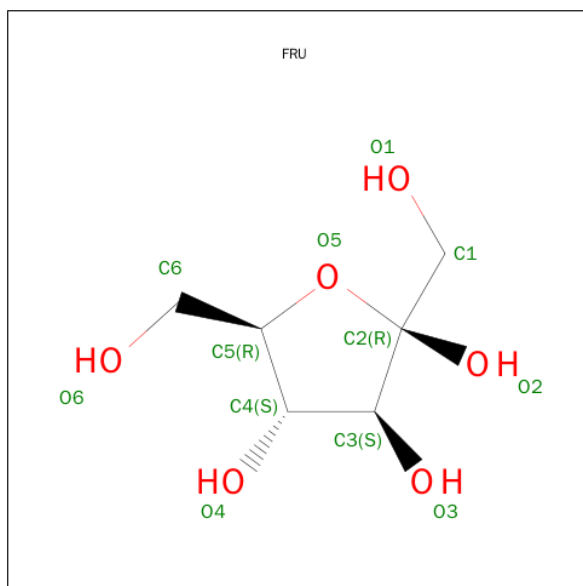
- Molecule 2 is a protein called Thiocyanate hydrolase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	153	Total	C	N	O	S	0	0	0
			1241	784	224	227	6			
2	E	151	Total	C	N	O	S	0	0	0
			1226	775	221	224	6			
2	H	156	Total	C	N	O	S	0	0	0
			1262	796	228	232	6			
2	K	152	Total	C	N	O	S	0	0	0
			1232	778	222	226	6			

- Molecule 3 is a protein called Thiocyanate hydrolase subunit gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	217	Total	C	N	O	S	0	0	0
			1725	1098	304	315	8			
3	F	218	Total	C	N	O	S	0	0	0
			1735	1104	307	316	8			
3	I	217	Total	C	N	O	S	0	0	0
			1725	1098	304	315	8			
3	L	216	Total	C	N	O	S	0	0	0
			1716	1093	303	312	8			

- Molecule 4 is SUGAR (FRUCTOSE) (three-letter code: FRU) (formula: $C_6H_{12}O_6$).

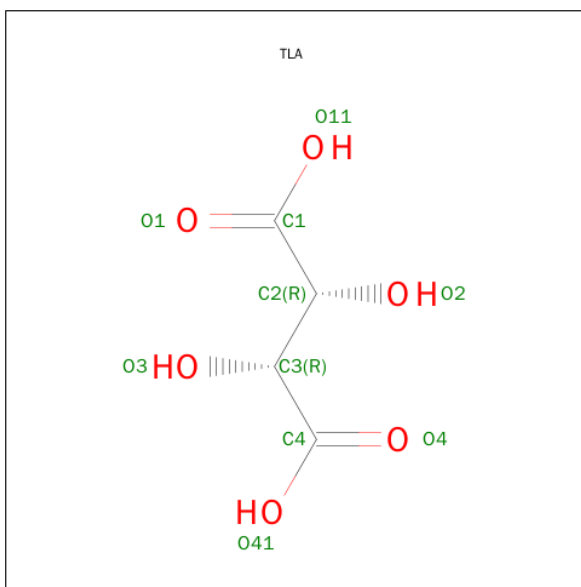


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			12	6	6		
4	L	1	Total	C	O	0	0
			12	6	6		
4	L	1	Total	C	O	0	0
			12	6	6		

- Molecule 5 is COBALT (III) ION (three-letter code: 3CO) (formula: Co).

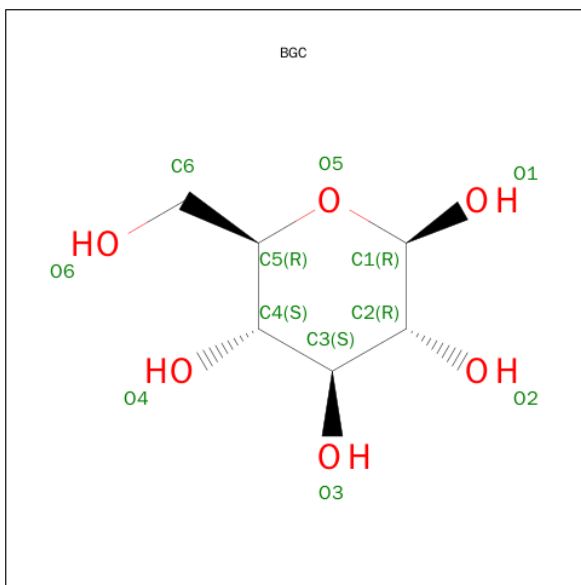
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	I	1	Total	Co	0	0
			1	1		
5	L	1	Total	Co	0	0
			1	1		
5	C	1	Total	Co	0	0
			1	1		
5	F	1	Total	Co	0	0
			1	1		

- Molecule 6 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: $C_4H_6O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			10	4	6		
6	F	1	Total	C	O	0	0
			10	4	6		
6	I	1	Total	C	O	0	0
			10	4	6		
6	L	1	Total	C	O	0	0
			10	4	6		

- Molecule 7 is SUGAR (BETA-D-GLUCOSE) (three-letter code: BGC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	F	1	Total	C	O	0	0
			12	6	6		

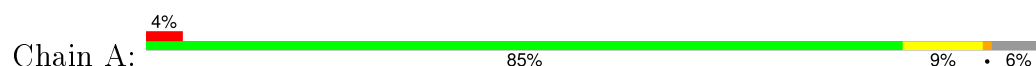
- Molecule 8 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	214	Total	O		0	0
			214	214			
8	B	265	Total	O		0	0
			265	265			
8	C	399	Total	O		0	0
			399	399			
8	D	216	Total	O		0	0
			216	216			
8	E	253	Total	O		0	0
			253	253			
8	F	352	Total	O		0	0
			352	352			
8	G	205	Total	O		0	0
			205	205			
8	H	260	Total	O		0	0
			260	260			
8	I	402	Total	O		0	0
			402	402			
8	J	206	Total	O		0	0
			206	206			
8	K	253	Total	O		0	0
			253	253			
8	L	377	Total	O		0	0
			377	377			

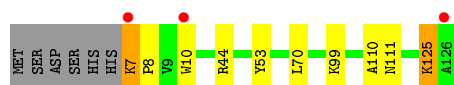
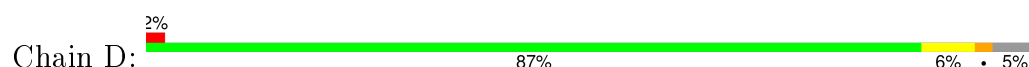
3 Residue-property plots [i](#)

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

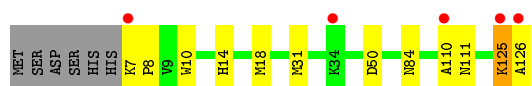
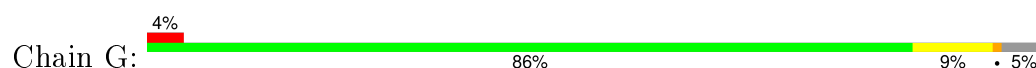
- Molecule 1: Thiocyanate hydrolase subunit alpha



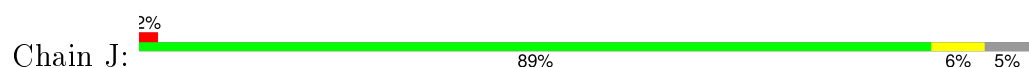
- Molecule 1: Thiocyanate hydrolase subunit alpha



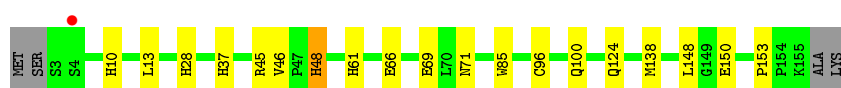
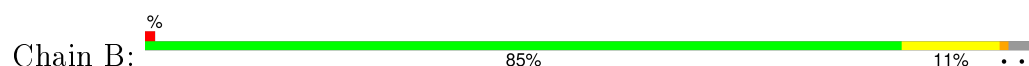
- Molecule 1: Thiocyanate hydrolase subunit alpha



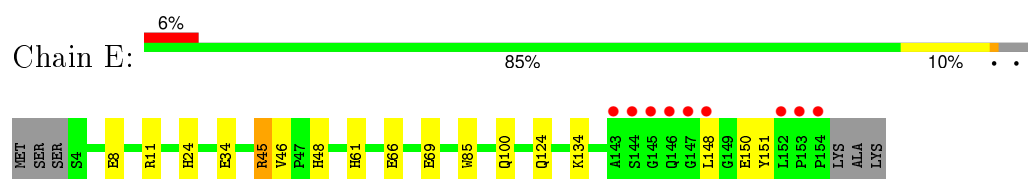
- Molecule 1: Thiocyanate hydrolase subunit alpha



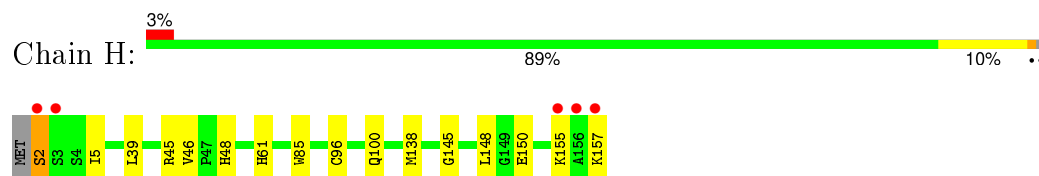
- Molecule 2: Thiocyanate hydrolase subunit beta



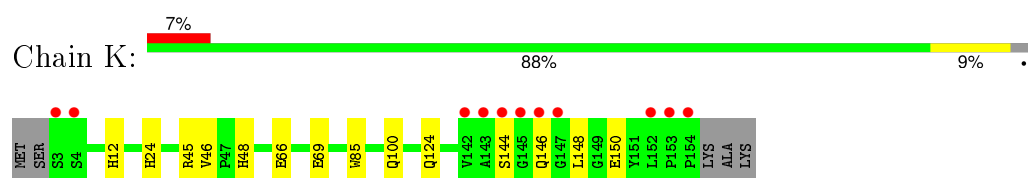
- Molecule 2: Thiocyanate hydrolase subunit beta



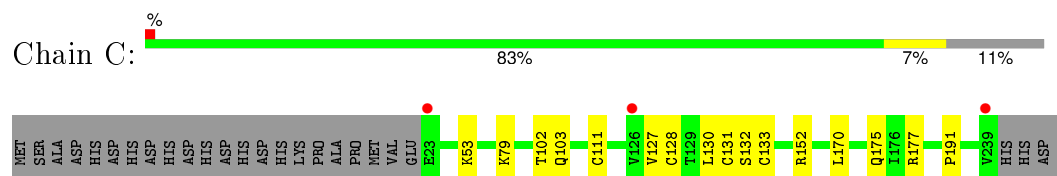
- Molecule 2: Thiocyanate hydrolase subunit beta



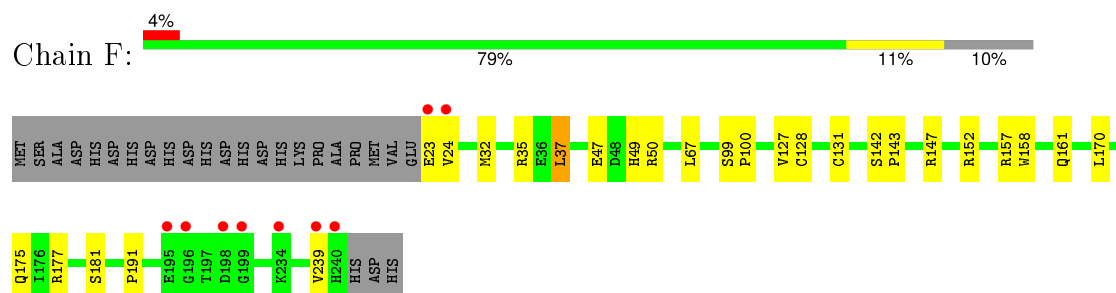
- Molecule 2: Thiocyanate hydrolase subunit beta



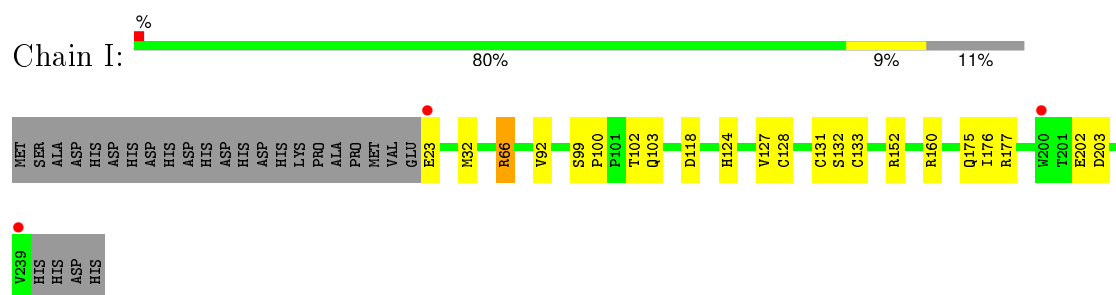
- Molecule 3: Thiocyanate hydrolase subunit gamma



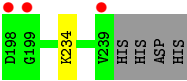
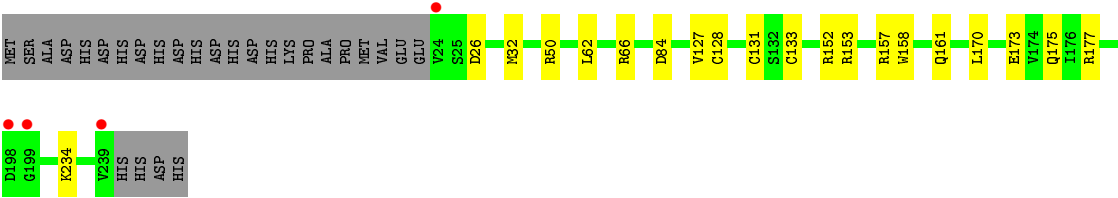
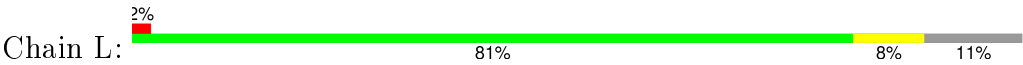
- Molecule 3: Thiocyanate hydrolase subunit gamma



- Molecule 3: Thiocyanate hydrolase subunit gamma



- Molecule 3: Thiocyanate hydrolase subunit gamma



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.01Å 170.76Å 175.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.76 – 1.78 40.76 – 1.78	Depositor EDS
% Data completeness (in resolution range)	94.3 (40.76-1.78) 97.3 (40.76-1.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.05 (at 1.78Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.150 , 0.159 0.153 , 0.164	Depositor DCC
R_{free} test set	16106 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 58.6	EDS
Estimated twinning fraction	0.007 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 319040 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19243	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.11% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CSD, 3CO, BGC, TLA, FRU

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	0.34	0/991	0.67	1/1342 (0.1%)
1	D	0.34	0/1000	0.64	1/1354 (0.1%)
1	G	0.34	0/1000	0.64	0/1354
1	J	0.35	0/1000	0.64	1/1354 (0.1%)
2	B	0.32	0/1273	0.63	0/1731
2	E	0.31	0/1258	0.62	0/1712
2	H	0.32	0/1294	0.64	0/1757
2	K	0.31	0/1264	0.63	0/1720
3	C	0.31	0/1753	0.65	0/2394
3	F	0.32	0/1764	0.67	0/2409
3	I	0.32	0/1753	0.66	0/2394
3	L	0.32	0/1744	0.66	0/2382
All	All	0.32	0/16094	0.65	3/21903 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	LEU	N-CA-C	-5.40	96.43	111.00
1	D	70	LEU	N-CA-C	-5.08	97.29	111.00
1	J	70	LEU	N-CA-C	-5.01	97.47	111.00

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	965	0	914	18	0
1	D	974	0	926	11	0
1	G	974	0	926	14	0
1	J	974	0	926	8	0
2	B	1241	0	1211	28	0
2	E	1226	0	1193	18	0
2	H	1262	0	1234	14	0
2	K	1232	0	1198	13	0
3	C	1725	0	1735	11	0
3	F	1735	0	1742	23	0
3	I	1725	0	1735	14	0
3	L	1716	0	1729	16	0
4	B	12	0	12	5	0
4	L	24	0	24	2	0
5	C	1	0	0	0	0
5	F	1	0	0	0	0
5	I	1	0	0	0	0
5	L	1	0	0	0	0
6	C	10	0	4	0	0
6	F	10	0	4	0	0
6	I	10	0	4	0	0
6	L	10	0	4	0	0
7	F	12	0	12	0	0
8	A	214	0	0	6	0
8	B	265	0	0	3	0
8	C	399	0	0	3	0
8	D	216	0	0	5	0
8	E	253	0	0	10	0
8	F	352	0	0	4	0
8	G	205	0	0	5	0
8	H	260	0	0	5	0
8	I	402	0	0	2	0
8	J	206	0	0	4	0
8	K	253	0	0	5	0
8	L	377	0	0	5	0
All	All	19243	0	15533	162	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 5.

All (162) 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)
2:E:148:LEU:CD1	3:F:32:MET:HG3	1.87	1.04
1:A:26:GLN:HE21	1:A:26:GLN:H	1.11	0.98
8:G:3339:HOH:O	2:H:46:VAL:HG23	1.62	0.98
2:B:138:MET:HE3	2:B:138:MET:HA	1.45	0.96
2:E:148:LEU:HD11	3:F:32:MET:HG3	1.44	0.95
1:G:125:LYS:HD3	1:G:126:ALA:N	1.84	0.93
3:L:234:LYS:HE2	8:L:3196:HOH:O	1.74	0.87
3:F:142:SER:OG	3:F:147:ARG:HD2	1.77	0.84
2:H:2:SER:HA	8:H:3352:HOH:O	1.81	0.79
2:E:45:ARG:HH21	2:E:100:GLN:HE22	1.32	0.77
2:E:148:LEU:HD13	3:F:32:MET:HG3	1.66	0.76
2:E:24:HIS:ND1	2:K:24:HIS:HD2	1.86	0.73
2:B:13:LEU:HB2	4:B:6002:FRU:H12	1.69	0.73
8:D:3249:HOH:O	1:G:7:LYS:HB2	1.89	0.71
3:F:175:GLN:HE22	3:F:177:ARG:HH11	1.37	0.71
1:G:125:LYS:HD3	1:G:125:LYS:C	2.07	0.70
1:A:26:GLN:NE2	1:A:26:GLN:H	1.85	0.69
3:I:175:GLN:HE22	3:I:177:ARG:HH21	1.39	0.69
2:B:28:HIS:HD2	3:L:153:ARG:HH12	1.41	0.68
2:B:138:MET:HE1	2:B:148:LEU:CD1	2.24	0.68
1:D:7:LYS:N	1:D:7:LYS:HD2	2.08	0.68
3:C:175:GLN:HE22	3:C:177:ARG:HH11	1.40	0.68
4:B:6002:FRU:O1	3:F:157:ARG:NH1	2.22	0.67
1:D:125:LYS:HE2	8:D:1430:HOH:O	1.94	0.66
2:B:10:HIS:ND1	4:B:6002:FRU:H11	2.11	0.66
3:L:175:GLN:HE22	3:L:177:ARG:HH21	1.44	0.66
2:B:138:MET:CE	2:B:148:LEU:CD1	2.75	0.65
2:B:28:HIS:CD2	3:L:153:ARG:HH12	2.13	0.65
1:A:26:GLN:HE21	1:A:26:GLN:N	1.90	0.65
2:B:150:GLU:HG2	8:B:3251:HOH:O	1.97	0.65
3:I:66:ARG:NH2	3:I:203:ASP:OD1	2.30	0.64
1:A:10:TRP:HZ3	8:K:3182:HOH:O	1.81	0.64
1:A:26:GLN:HE22	3:C:111:CYS:HB2	1.64	0.63
1:D:10:TRP:HZ3	8:E:3201:HOH:O	1.81	0.63
8:E:1521:HOH:O	3:F:49:HIS:HD2	1.83	0.62
2:B:153:PRO:HG2	8:B:2426:HOH:O	1.99	0.62
8:E:1116:HOH:O	3:F:239:VAL:HG12	1.98	0.62
2:B:138:MET:CE	2:B:138:MET:HA	2.26	0.62
1:D:8:PRO:HG3	1:G:10:TRP:CH2	2.36	0.60
2:B:138:MET:HE3	2:B:138:MET:CA	2.25	0.60
1:A:125:LYS:HB2	8:A:2537:HOH:O	2.02	0.60
1:J:125:LYS:HG3	8:J:2432:HOH:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:31:MET:HE2	8:G:3130:HOH:O	2.03	0.59
3:F:47:GLU:OE1	3:F:50:ARG:NH2	2.36	0.59
2:B:37:HIS:HE1	2:E:8:GLU:OE2	1.85	0.59
2:E:34:GLU:HG3	8:E:3257:HOH:O	2.03	0.58
2:K:45:ARG:HH11	2:K:100:GLN:HE22	1.52	0.58
2:H:46:VAL:CG1	2:H:96:CYS:SG	2.92	0.58
2:B:71:ASN:HD21	3:C:53:LYS:NZ	2.01	0.57
2:H:45:ARG:HH11	2:H:100:GLN:HE22	1.49	0.57
1:D:53:TYR:HA	2:E:46:VAL:HG11	1.86	0.57
1:D:8:PRO:HG3	1:G:10:TRP:CZ3	2.40	0.57
3:I:66:ARG:NE	3:I:202:GLU:OE2	2.29	0.56
3:F:143:PRO:O	3:F:147:ARG:HD3	2.06	0.56
1:J:31:MET:HE2	8:J:2594:HOH:O	2.05	0.55
2:B:46:VAL:CG1	2:B:96:CYS:SG	2.94	0.55
3:I:160:ARG:HG2	3:I:176:ILE:HD12	1.87	0.55
4:B:6002:FRU:HO1	3:F:157:ARG:HH12	1.48	0.55
2:K:148:LEU:HG	3:L:32:MET:HG3	1.89	0.54
2:K:146:GLN:NE2	8:K:3293:HOH:O	2.39	0.54
2:K:46:VAL:HG22	8:K:3223:HOH:O	2.06	0.54
3:F:127:VAL:HG22	3:F:128:CYS:N	2.23	0.54
2:K:46:VAL:HG23	2:K:46:VAL:O	2.08	0.54
2:B:45:ARG:HH11	2:B:100:GLN:HE22	1.56	0.53
3:I:102:THR:O	3:I:103:GLN:HB2	2.07	0.53
1:G:7:LYS:N	1:G:8:PRO:CD	2.72	0.52
3:L:127:VAL:HG22	3:L:128:CYS:N	2.25	0.52
3:C:102:THR:O	3:C:103:GLN:HB2	2.10	0.52
3:I:127:VAL:HG22	3:I:128:CYS:N	2.25	0.52
1:J:7:LYS:HE3	8:J:3266:HOH:O	2.10	0.51
1:A:10:TRP:CZ3	1:J:8:PRO:HG2	2.44	0.51
4:L:6003:FRU:H62	8:L:2670:HOH:O	2.10	0.51
1:G:125:LYS:HD3	1:G:126:ALA:H	1.72	0.51
3:I:118:ASP:OD1	3:I:124:HIS:HD2	1.93	0.51
3:C:127:VAL:HG22	3:C:128:CYS:N	2.25	0.51
1:G:14:HIS:HE1	8:G:900:HOH:O	1.92	0.51
2:E:150:GLU:HG2	8:E:2397:HOH:O	2.10	0.51
1:D:10:TRP:HH2	8:E:2814:HOH:O	1.94	0.51
2:B:46:VAL:HG12	2:B:96:CYS:SG	2.50	0.51
2:E:148:LEU:HD12	2:E:151:TYR:HD2	1.74	0.51
3:L:84:ASP:OD2	4:L:6001:FRU:O1	2.21	0.51
3:I:23:GLU:O	3:I:23:GLU:HG3	2.10	0.50
1:G:7:LYS:N	1:G:8:PRO:HD3	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:150:GLU:HG2	8:K:1250:HOH:O	2.11	0.50
3:C:79:LYS:HE2	8:C:3344:HOH:O	2.11	0.50
2:H:145:GLY:HA2	2:H:155:LYS:HE3	1.93	0.50
2:B:138:MET:CA	2:B:138:MET:CE	2.87	0.49
2:B:61:HIS:HE1	8:C:1187:HOH:O	1.95	0.49
1:A:44:ARG:NE	8:A:3389:HOH:O	2.45	0.49
2:H:45:ARG:NH1	2:H:100:GLN:HE22	2.11	0.49
2:E:61:HIS:HE1	8:F:1164:HOH:O	1.94	0.49
1:G:111:ASN:O	3:I:175:GLN:HG3	2.13	0.48
1:A:10:TRP:HH2	8:K:3191:HOH:O	1.95	0.48
1:A:44:ARG:CZ	8:A:3389:HOH:O	2.61	0.48
3:L:131:CSD:O	3:L:152:ARG:HD2	2.13	0.48
3:I:23:GLU:O	3:I:23:GLU:CG	2.60	0.48
1:A:84:ASN:ND2	2:K:124:GLN:HE21	2.10	0.48
1:A:34:LYS:HD3	8:A:3168:HOH:O	2.12	0.48
2:E:124:GLN:HE21	1:G:84:ASN:ND2	2.11	0.48
3:F:131:CSD:O	3:F:152:ARG:HD2	2.14	0.48
3:F:50:ARG:HG3	8:F:986:HOH:O	2.14	0.48
1:A:10:TRP:CH2	1:J:8:PRO:HG2	2.49	0.47
2:H:157:LYS:HG3	2:H:157:LYS:OXT	2.14	0.47
8:I:1162:HOH:O	3:L:161:GLN:HG2	2.14	0.47
2:B:138:MET:HE2	2:B:148:LEU:CD1	2.44	0.47
3:C:131:CSD:O	3:C:152:ARG:HD2	2.15	0.47
2:H:2:SER:HB3	2:H:5:ILE:HD12	1.97	0.47
2:H:150:GLU:HG2	8:H:2405:HOH:O	2.14	0.47
8:E:174:HOH:O	2:K:24:HIS:HE1	1.97	0.47
8:C:3224:HOH:O	3:F:161:GLN:HG2	2.15	0.47
3:I:131:CSD:O	3:I:152:ARG:HD2	2.15	0.47
2:B:10:HIS:HA	4:B:6002:FRU:H11	1.97	0.46
8:H:396:HOH:O	2:K:12:HIS:HE1	1.98	0.46
3:F:67:LEU:HD23	3:F:67:LEU:C	2.35	0.46
2:B:124:GLN:HE21	1:J:84:ASN:ND2	2.13	0.46
2:H:61:HIS:HE1	8:I:862:HOH:O	1.98	0.46
1:D:111:ASN:O	3:F:175:GLN:HG3	2.16	0.46
1:J:111:ASN:O	3:L:175:GLN:HG3	2.15	0.46
2:K:144:SER:OG	2:K:146:GLN:HG3	2.15	0.46
8:B:3190:HOH:O	2:E:11:ARG:HD3	2.15	0.46
3:F:175:GLN:NE2	3:F:177:ARG:HH11	2.11	0.46
1:A:111:ASN:O	3:C:175:GLN:HG3	2.15	0.46
1:D:99:LYS:HE2	8:D:3345:HOH:O	2.17	0.45
1:A:10:TRP:CZ3	1:J:8:PRO:CG	3.00	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:PRO:HG3	8:J:3264:HOH:O	2.15	0.45
2:H:46:VAL:HG12	2:H:96:CYS:SG	2.57	0.45
1:D:44:ARG:NH2	8:D:3353:HOH:O	2.49	0.45
2:B:37:HIS:HD2	3:L:26:ASP:OD2	1.99	0.45
2:K:45:ARG:NH1	2:K:100:GLN:HE22	2.14	0.45
2:E:34:GLU:HG3	8:E:844:HOH:O	2.15	0.44
3:L:62:LEU:O	3:L:66:ARG:HG3	2.16	0.44
2:E:66:GLU:HB2	2:E:69:GLU:HG3	1.99	0.44
2:E:134:LYS:HE3	3:F:37:LEU:HD13	2.00	0.44
2:B:45:ARG:NH1	2:B:100:GLN:HE22	2.17	0.43
3:I:132:SER:O	3:I:133:CSD:C	2.67	0.43
2:H:138:MET:SD	2:H:148:LEU:HD11	2.58	0.42
2:H:2:SER:CA	8:H:3352:HOH:O	2.55	0.42
3:C:132:SER:O	3:C:133:CSD:C	2.67	0.42
1:A:99:LYS:NZ	8:A:3380:HOH:O	2.52	0.42
3:F:99:SER:N	3:F:100:PRO:CD	2.83	0.42
3:I:99:SER:N	3:I:100:PRO:CD	2.82	0.42
3:L:173:GLU:HG2	8:L:2165:HOH:O	2.19	0.42
2:B:138:MET:CE	2:B:148:LEU:HD11	2.48	0.42
1:A:79:ASP:HA	8:A:341:HOH:O	2.20	0.42
2:B:71:ASN:HD21	3:C:53:LYS:HZ2	1.66	0.41
3:F:35:ARG:HB3	8:F:3348:HOH:O	2.20	0.41
1:D:110:ALA:HB2	8:D:1428:HOH:O	2.20	0.41
3:L:50:ARG:HG3	8:L:1173:HOH:O	2.20	0.41
1:G:50:ASP:CG	8:G:3339:HOH:O	2.59	0.41
3:I:66:ARG:HG2	3:I:92:VAL:HG12	2.02	0.41
2:E:61:HIS:HD2	8:E:2399:HOH:O	2.04	0.41
3:F:24:VAL:HG22	8:F:3269:HOH:O	2.20	0.41
1:G:110:ALA:HA	8:G:2549:HOH:O	2.20	0.41
2:B:138:MET:CE	2:B:148:LEU:HD13	2.50	0.41
2:B:48:HIS:HB2	3:C:130:LEU:O	2.20	0.41
2:K:66:GLU:HB2	2:K:69:GLU:HG3	2.02	0.41
3:F:157:ARG:HD3	3:F:158:TRP:CH2	2.56	0.41
2:H:61:HIS:CD2	8:H:1339:HOH:O	2.74	0.40
2:B:66:GLU:HB2	2:B:69:GLU:HG3	2.03	0.40
2:E:61:HIS:CD2	8:E:2399:HOH:O	2.73	0.40
3:L:157:ARG:HD3	3:L:158:TRP:CH2	2.57	0.40
3:L:133:CSD:OD2	8:L:245:HOH:O	2.17	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	117/126 (93%)	113 (97%)	4 (3%)	0	100	100
1	D	118/126 (94%)	114 (97%)	4 (3%)	0	100	100
1	G	118/126 (94%)	115 (98%)	3 (2%)	0	100	100
1	J	118/126 (94%)	115 (98%)	3 (2%)	0	100	100
2	B	151/157 (96%)	148 (98%)	3 (2%)	0	100	100
2	E	149/157 (95%)	146 (98%)	3 (2%)	0	100	100
2	H	154/157 (98%)	150 (97%)	4 (3%)	0	100	100
2	K	150/157 (96%)	147 (98%)	3 (2%)	0	100	100
3	C	213/243 (88%)	206 (97%)	7 (3%)	0	100	100
3	F	214/243 (88%)	206 (96%)	8 (4%)	0	100	100
3	I	213/243 (88%)	205 (96%)	8 (4%)	0	100	100
3	L	212/243 (87%)	204 (96%)	8 (4%)	0	100	100
All	All	1927/2104 (92%)	1869 (97%)	58 (3%)	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	101/108 (94%)	99 (98%)	2 (2%)	63	47
1	D	102/108 (94%)	100 (98%)	2 (2%)	63	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	102/108 (94%)	100 (98%)	2 (2%)	63	47
1	J	102/108 (94%)	101 (99%)	1 (1%)	82	75
2	B	131/134 (98%)	129 (98%)	2 (2%)	72	60
2	E	129/134 (96%)	126 (98%)	3 (2%)	58	40
2	H	133/134 (99%)	129 (97%)	4 (3%)	48	29
2	K	130/134 (97%)	128 (98%)	2 (2%)	72	60
3	C	188/212 (89%)	186 (99%)	2 (1%)	80	72
3	F	189/212 (89%)	184 (97%)	5 (3%)	54	35
3	I	188/212 (89%)	186 (99%)	2 (1%)	80	72
3	L	187/212 (88%)	186 (100%)	1 (0%)	92	89
All	All	1682/1816 (93%)	1654 (98%)	28 (2%)	68	54

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	MET
1	A	26	GLN
2	B	48	HIS
2	B	85	TRP
3	C	170	LEU
3	C	191	PRO
1	D	7	LYS
1	D	125	LYS
2	E	45	ARG
2	E	48	HIS
2	E	85	TRP
3	F	23	GLU
3	F	37	LEU
3	F	170	LEU
3	F	181	SER
3	F	191	PRO
1	G	18	MET
1	G	125	LYS
2	H	2	SER
2	H	39	LEU
2	H	48	HIS
2	H	85	TRP
3	I	32	MET

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Mol	Chain	Res	Type
3	I	66	ARG
1	J	99	LYS
2	K	48	HIS
2	K	85	TRP
3	L	170	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	84	ASN
1	A	111	ASN
2	B	28	HIS
2	B	37	HIS
2	B	61	HIS
2	B	71	ASN
2	B	100	GLN
3	C	103	GLN
3	C	169	GLN
3	C	175	GLN
2	E	61	HIS
2	E	94	GLN
2	E	100	GLN
3	F	49	HIS
3	F	103	GLN
3	F	175	GLN
1	G	14	HIS
1	G	84	ASN
2	H	61	HIS
2	H	100	GLN
2	H	124	GLN
2	H	146	GLN
3	I	103	GLN
3	I	112	ASN
3	I	124	HIS
3	I	169	GLN
3	I	175	GLN
1	J	84	ASN
2	K	12	HIS
2	K	24	HIS
2	K	100	GLN
2	K	146	GLN

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Mol	Chain	Res	Type
3	L	57	HIS
3	L	112	ASN
3	L	175	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	CSD	C	131	3,5	3,7,8	0.71	0	3,8,10	1.71	1 (33%)
3	CSD	C	133	3,5	3,7,8	0.83	0	3,8,10	2.11	2 (66%)
3	CSD	F	131	3,5	3,7,8	0.71	0	3,8,10	1.70	1 (33%)
3	CSD	F	133	3,5	3,7,8	0.82	0	3,8,10	2.09	2 (66%)
3	CSD	I	131	3,5	3,7,8	0.73	0	3,8,10	1.64	1 (33%)
3	CSD	I	133	3,5	3,7,8	0.75	0	3,8,10	2.15	2 (66%)
3	CSD	L	131	3,5	3,7,8	0.72	0	3,8,10	1.57	0
3	CSD	L	133	3,5	3,7,8	0.74	0	3,8,10	2.10	2 (66%)

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	CSD	C	131	3,5	-	0/2/6/8	0/0/0/0
3	CSD	C	133	3,5	-	0/2/6/8	0/0/0/0
3	CSD	F	131	3,5	-	0/2/6/8	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CSD	F	133	3,5	-	0/2/6/8	0/0/0/0
3	CSD	I	131	3,5	-	0/2/6/8	0/0/0/0
3	CSD	I	133	3,5	-	0/2/6/8	0/0/0/0
3	CSD	L	131	3,5	-	0/2/6/8	0/0/0/0
3	CSD	L	133	3,5	-	0/2/6/8	0/0/0/0

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	131	CSD	OD1-SG-CB	2.14	108.97	105.40
3	L	133	CSD	OD1-SG-CB	2.17	109.03	105.40
3	F	133	CSD	OD1-SG-CB	2.18	109.03	105.40
3	L	133	CSD	CB-CA-C	2.19	117.47	111.46
3	F	131	CSD	OD1-SG-CB	2.20	109.07	105.40
3	C	133	CSD	OD1-SG-CB	2.20	109.07	105.40
3	I	133	CSD	OD1-SG-CB	2.25	109.16	105.40
3	C	131	CSD	OD1-SG-CB	2.28	109.20	105.40
3	F	133	CSD	CB-CA-C	2.28	117.72	111.46
3	C	133	CSD	CB-CA-C	2.32	117.83	111.46
3	I	133	CSD	CB-CA-C	2.37	117.94	111.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	131	CSD	1	0
3	C	133	CSD	1	0
3	F	131	CSD	1	0
3	I	131	CSD	1	0
3	I	133	CSD	1	0
3	L	131	CSD	1	0
3	L	133	CSD	1	0

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	FRU	B	6002	-	11,12,12	1.70	2 (18%)	10,18,18	1.57	1 (10%)
6	TLA	C	4001	-	3,9,9	0.82	0	6,12,12	1.13	1 (16%)
6	TLA	F	4002	-	3,9,9	0.78	0	6,12,12	1.11	1 (16%)
7	BGC	F	5001	-	12,12,12	0.36	0	17,17,17	0.38	0
6	TLA	I	4003	-	3,9,9	0.81	0	6,12,12	1.17	1 (16%)
6	TLA	L	4004	-	3,9,9	0.80	0	6,12,12	1.17	1 (16%)
4	FRU	L	6001	-	11,12,12	1.66	3 (27%)	10,18,18	2.33	3 (30%)
4	FRU	L	6003	-	11,12,12	1.73	2 (18%)	10,18,18	1.67	2 (20%)

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	FRU	B	6002	-	-	0/5/24/24	0/1/1/1
6	TLA	C	4001	-	-	0/4/12/12	0/0/0/0
6	TLA	F	4002	-	-	0/4/12/12	0/0/0/0
7	BGC	F	5001	-	-	0/2/22/22	0/1/1/1
6	TLA	I	4003	-	-	0/4/12/12	0/0/0/0
6	TLA	L	4004	-	-	0/4/12/12	0/0/0/0
4	FRU	L	6001	-	-	0/5/24/24	0/1/1/1
4	FRU	L	6003	-	-	0/5/24/24	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	6001	FRU	O3-C3	2.06	1.47	1.42
4	B	6002	FRU	O5-C5	2.15	1.48	1.43
4	L	6001	FRU	O2-C2	2.97	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	6003	FRU	O5-C5	3.09	1.50	1.43
4	L	6001	FRU	C1-C2	3.46	1.58	1.52
4	L	6003	FRU	C1-C2	4.40	1.59	1.52
4	B	6002	FRU	C1-C2	4.55	1.60	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	6003	FRU	C6-C5-C4	-2.78	108.51	115.08
6	C	4001	TLA	C4-C3-C2	-2.30	108.64	113.35
6	L	4004	TLA	C4-C3-C2	-2.29	108.66	113.35
4	L	6001	FRU	O6-C6-C5	-2.24	103.94	111.33
6	I	4003	TLA	C4-C3-C2	-2.19	108.85	113.35
6	F	4002	TLA	C4-C3-C2	-2.15	108.95	113.35
4	L	6001	FRU	C6-C5-C4	2.92	122.00	115.08
4	L	6003	FRU	O2-C2-O5	2.93	115.42	109.37
4	B	6002	FRU	O2-C2-O5	3.88	117.37	109.37
4	L	6001	FRU	O2-C2-O5	5.70	121.13	109.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	6002	FRU	5	0
4	L	6001	FRU	1	0
4	L	6003	FRU	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	119/126 (94%)	0.04	5 (4%) 40 38	11, 17, 32, 39	0
1	D	120/126 (95%)	-0.16	3 (2%) 61 59	11, 15, 31, 45	0
1	G	120/126 (95%)	-0.02	5 (4%) 40 38	11, 17, 31, 48	0
1	J	120/126 (95%)	-0.22	2 (1%) 73 72	10, 14, 28, 48	0
2	B	153/157 (97%)	-0.24	1 (0%) 89 89	10, 13, 26, 44	0
2	E	151/157 (96%)	-0.03	9 (5%) 25 23	10, 15, 32, 41	0
2	H	156/157 (99%)	-0.10	5 (3%) 51 49	10, 13, 26, 70	0
2	K	152/157 (96%)	-0.05	11 (7%) 18 17	10, 14, 29, 40	0
3	C	215/243 (88%)	-0.11	3 (1%) 78 77	10, 15, 26, 45	0
3	F	216/243 (88%)	0.11	9 (4%) 40 38	10, 18, 34, 54	0
3	I	215/243 (88%)	-0.13	3 (1%) 78 77	10, 14, 24, 46	0
3	L	214/243 (88%)	-0.12	4 (1%) 70 69	10, 14, 26, 36	0
All	All	1951/2104 (92%)	-0.08	60 (3%) 52 50	10, 15, 29, 70	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	157	LYS	8.2
3	F	240	HIS	7.6
2	H	156	ALA	7.1
2	H	2	SER	5.6
2	H	3	SER	5.5
3	I	23	GLU	4.9
3	F	23	GLU	4.7
3	I	239	VAL	4.2
3	C	239	VAL	3.9
3	F	239	VAL	3.7
1	A	126	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
3	L	239	VAL	3.7
2	E	147	GLY	3.7
2	E	148	LEU	3.5
2	E	153	PRO	3.5
2	K	146	GLN	3.5
2	K	152	LEU	3.4
2	K	145	GLY	3.3
2	E	145	GLY	3.2
2	K	3	SER	3.2
1	A	9	VAL	3.2
2	E	144	SER	3.1
3	L	198	ASP	3.0
3	F	198	ASP	3.0
2	K	143	ALA	3.0
2	K	154	PRO	2.9
3	C	23	GLU	2.9
1	A	34	LYS	2.9
2	E	154	PRO	2.9
1	A	125	LYS	2.7
1	G	126	ALA	2.7
2	K	144	SER	2.7
2	E	146	GLN	2.7
1	D	7	LYS	2.7
2	K	153	PRO	2.6
1	D	126	ALA	2.6
3	F	196	GLY	2.5
2	E	152	LEU	2.5
1	G	110	ALA	2.4
1	J	126	ALA	2.4
1	G	34	LYS	2.4
3	F	199	GLY	2.4
2	E	143	ALA	2.4
2	H	155	LYS	2.4
2	K	4	SER	2.3
3	L	24	VAL	2.3
2	K	142	VAL	2.3
1	G	7	LYS	2.2
2	K	147	GLY	2.2
3	L	199	GLY	2.2
1	D	10	TRP	2.2
3	C	126	VAL	2.2
3	I	200	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	8	PRO	2.1
3	F	195	GLU	2.1
1	G	125	LYS	2.1
1	J	7	LYS	2.0
3	F	234	LYS	2.0
2	B	4	SER	2.0
3	F	24	VAL	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
3	CSD	L	131	8/9	0.99	0.06	-	10,11,12,12	0
3	CSD	I	133	8/9	0.97	0.10	-	11,12,15,15	1
3	CSD	F	133	8/9	0.98	0.07	-	12,14,15,15	1
3	CSD	C	131	8/9	0.99	0.08	-	10,12,13,15	0
3	CSD	I	131	8/9	0.99	0.08	-	11,12,13,14	0
3	CSD	L	133	8/9	0.98	0.07	-	10,11,14,14	1
3	CSD	C	133	8/9	0.98	0.10	-	11,13,15,16	1
3	CSD	F	131	8/9	0.99	0.06	-	12,12,13,14	0

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(\AA^2)	Q<0.9
4	FRU	L	6003	12/12	0.56	0.42	16.93	24,41,46,51	0
4	FRU	B	6002	12/12	0.36	0.51	16.32	33,38,41,42	0
4	FRU	L	6001	12/12	0.58	0.39	14.08	43,51,59,60	0
7	BGC	F	5001	12/12	0.92	0.19	12.22	25,27,28,28	0
5	3CO	F	301	1/1	1.00	0.04	-	14,14,14,14	0
6	TLA	L	4004	10/10	0.95	0.12	-	16,18,23,24	0
6	TLA	F	4002	10/10	0.95	0.11	-	18,22,25,27	0
5	3CO	C	301	1/1	1.00	0.03	-	14,14,14,14	0
5	3CO	L	301	1/1	1.00	0.03	-	13,13,13,13	0
6	TLA	I	4003	10/10	0.94	0.15	-	20,21,25,26	0
6	TLA	C	4001	10/10	0.95	0.11	-	19,20,24,26	0
5	3CO	I	301	1/1	1.00	0.03	-	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.