



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2018 – 05:53 pm GMT

PDB ID : 2DD5
Title : Thiocyanate hydrolase (SCNase) from Thiobacillus thioparus native holo-enzyme
Authors : Arakawa, T.; Kawano, Y.; Kataoka, S.; Katayama, Y.; Kamiya, N.; Yohda, M.; Odaka, M.
Deposited on : 2006-01-19
Resolution : 2.00 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 : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

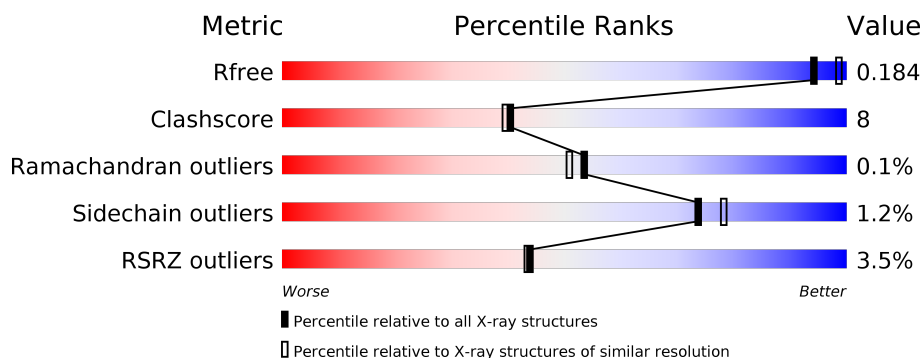
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	111664	7193 (2.00-2.00)
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

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>2%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>6%</div> </div> </div>
1	D	126	<div> <div>6%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>6%</div> </div> </div>
1	G	126	<div> <div>5%</div> <div> <div></div> <div>86%</div> <div>10%</div> <div>5%</div> </div> </div>
1	J	126	<div> <div>2%</div> <div> <div></div> <div>89%</div> <div>6%</div> <div>5%</div> </div> </div>
2	B	157	<div> <div>7%</div> <div> <div></div> <div>76%</div> <div>20%</div> <div>• •</div> </div> </div>
2	E	157	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>20%</div> <div>• •</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	H	157	<div><div></div><div>9%</div><div>82%</div><div>16%</div><div>••</div></div>
2	K	157	<div><div></div><div>%</div><div>80%</div><div>17%</div><div>•</div></div>
3	C	243	<div><div></div><div>3%</div><div>73%</div><div>16%</div><div>11%</div></div>
3	F	243	<div><div></div><div>2%</div><div>70%</div><div>18%</div><div>11%</div></div>
3	I	243	<div><div></div><div>2%</div><div>74%</div><div>15%</div><div>11%</div></div>
3	L	243	<div><div></div><div></div><div>79%</div><div>9%</div><div>11%</div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17962 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 alpha subunit.

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	119	Total	C	N	O	S	0	0	0
			965	614	160	187	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			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP O66187
D	1	MET	-	INITIATING METHIONINE	UNP O66187
G	1	MET	-	INITIATING METHIONINE	UNP O66187
J	1	MET	-	INITIATING METHIONINE	UNP O66187

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	152	Total	C	N	O	S	0	0	0
			1232	778	222	226	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			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	INITIATING METHIONINE	UNP O66186
E	1	MET	-	INITIATING METHIONINE	UNP O66186
H	1	MET	-	INITIATING METHIONINE	UNP O66186
K	1	MET	-	INITIATING METHIONINE	UNP O66186

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	217	Total	C	N	O	S	3	0	0
			1724	1098	304	314	8			
3	F	216	Total	C	N	O	S	3	0	0
			1715	1093	303	311	8			
3	I	217	Total	C	N	O	S	3	0	0
			1724	1098	304	314	8			
3	L	216	Total	C	N	O	S	3	0	0
			1715	1093	303	311	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	INITIATING METHIONINE	UNP O66188
C	131	CSD	CYS	MODIFIED RESIDUE	UNP O66188
C	133	CSO	CYS	MODIFIED RESIDUE	UNP O66188
F	1	MET	-	INITIATING METHIONINE	UNP O66188
F	131	CSD	CYS	MODIFIED RESIDUE	UNP O66188
F	133	CSO	CYS	MODIFIED RESIDUE	UNP O66188
I	1	MET	-	INITIATING METHIONINE	UNP O66188
I	131	CSD	CYS	MODIFIED RESIDUE	UNP O66188
I	133	CSO	CYS	MODIFIED RESIDUE	UNP O66188
L	1	MET	-	INITIATING METHIONINE	UNP O66188
L	131	CSD	CYS	MODIFIED RESIDUE	UNP O66188
L	133	CSO	CYS	MODIFIED RESIDUE	UNP O66188

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	J	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		

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

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

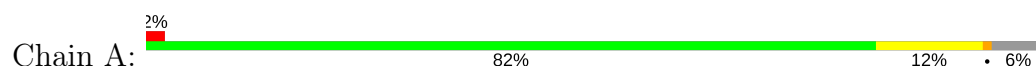
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	134	Total O 134 134	0	0
6	B	161	Total O 161 161	0	0
6	C	249	Total O 249 249	0	0
6	D	103	Total O 103 103	0	0
6	E	157	Total O 157 157	0	0
6	F	233	Total O 233 233	0	0
6	G	134	Total O 134 134	0	0
6	H	157	Total O 157 157	0	0
6	I	263	Total O 263 263	0	0
6	J	148	Total O 148 148	0	0
6	K	164	Total O 164 164	0	0
6	L	287	Total O 287 287	0	0

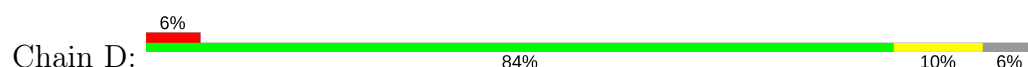
3 Residue-property plots [i](#)

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.

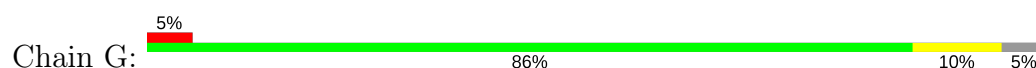
- Molecule 1: Thiocyanate hydrolase alpha subunit



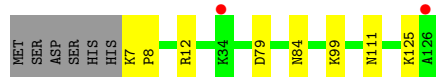
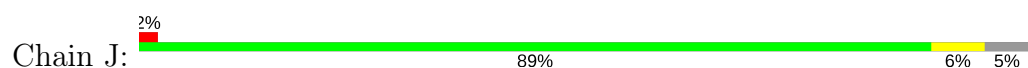
- Molecule 1: Thiocyanate hydrolase alpha subunit



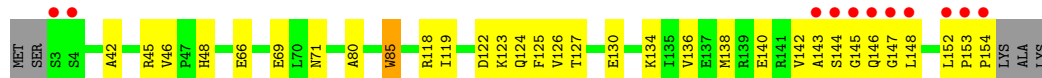
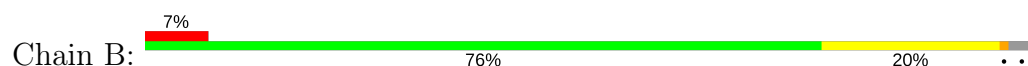
- Molecule 1: Thiocyanate hydrolase alpha subunit



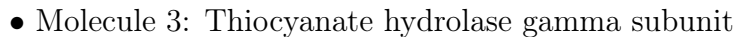
- Molecule 1: Thiocyanate hydrolase alpha subunit

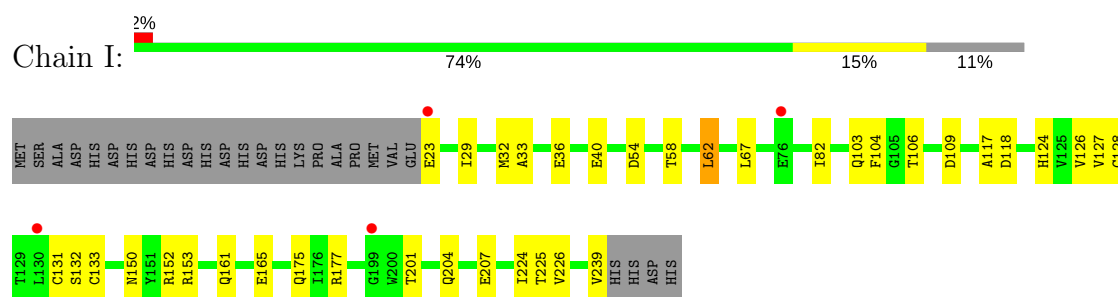


- Molecule 2: Thiocyanate hydrolase beta subunit

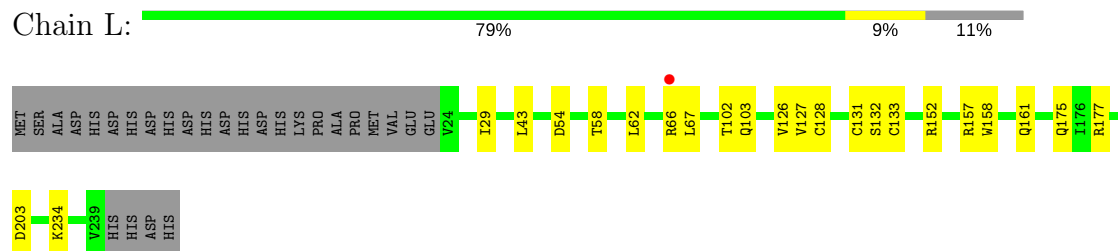


- Molecule 2: Thiocyanate hydrolase beta subunit





- Molecule 3: Thiocyanate hydrolase gamma subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	170.19Å 170.19Å 282.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.80 – 2.00 37.80 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.3 (37.80-2.00) 99.9 (37.80-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.99 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.185 , 0.205 0.183 , 0.184	Depositor DCC
R_{free} test set	14043 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17962	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.89% 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: CSD, 3CO, CSO, SO4

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	0.61	0/1342
1	D	0.32	0/991	0.59	0/1342
1	G	0.33	0/1000	0.60	0/1354
1	J	0.34	0/1000	0.61	0/1354
2	B	0.31	0/1264	0.58	0/1720
2	E	0.29	0/1258	0.56	0/1712
2	H	0.30	0/1294	0.58	0/1757
2	K	0.32	0/1264	0.59	0/1720
3	C	0.29	0/1752	0.60	0/2391
3	F	0.28	0/1743	0.59	0/2379
3	I	0.30	0/1752	0.61	0/2391
3	L	0.31	0/1743	0.62	0/2379
All	All	0.31	0/16052	0.59	0/21841

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	965	0	914	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	965	0	914	9	0
1	G	974	0	926	10	0
1	J	974	0	926	11	0
2	B	1232	0	1198	30	0
2	E	1226	0	1193	34	0
2	H	1262	0	1234	32	0
2	K	1232	0	1198	17	0
3	C	1724	0	1735	44	0
3	F	1715	0	1729	37	0
3	I	1724	0	1735	34	0
3	L	1715	0	1729	17	0
4	A	5	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
4	F	5	0	0	0	0
4	G	5	0	0	0	0
4	I	15	0	0	0	0
4	J	5	0	0	0	0
4	L	15	0	0	0	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	A	134	0	0	2	0
6	B	161	0	0	3	0
6	C	249	0	0	3	0
6	D	103	0	0	0	0
6	E	157	0	0	2	0
6	F	233	0	0	3	0
6	G	134	0	0	0	0
6	H	157	0	0	2	0
6	I	263	0	0	3	0
6	J	148	0	0	1	0
6	K	164	0	0	2	0
6	L	287	0	0	1	0
All	All	17962	0	15431	244	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 8.

All (244) 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:H:148:LEU:HB3	3:I:32:MET:SD	2.06	0.95
1:A:26:GLN:HE21	1:A:26:GLN:H	1.01	0.94
3:I:36:GLU:O	3:I:40:GLU:HG3	1.69	0.93
3:I:82:ILE:HD11	3:I:117:ALA:HB2	1.51	0.90
2:B:148:LEU:HD11	3:C:32:MET:HE1	1.56	0.87
2:B:148:LEU:HG	3:C:32:MET:SD	2.16	0.85
2:E:148:LEU:HD11	3:F:32:MET:HE1	1.60	0.82
3:I:175:GLN:HE22	3:I:177:ARG:HH21	1.28	0.80
1:A:84:ASN:HD22	1:J:7:LYS:NZ	1.80	0.79
1:A:26:GLN:N	1:A:26:GLN:HE21	1.79	0.78
1:A:26:GLN:NE2	1:A:26:GLN:H	1.81	0.78
3:C:237:ARG:HH21	3:C:237:ARG:CG	1.98	0.77
2:H:58:ILE:HD11	2:H:60:TYR:CE1	2.20	0.76
3:C:237:ARG:HB2	3:C:237:ARG:HH21	1.50	0.76
6:C:3152:HOH:O	3:F:161:GLN:HG2	1.88	0.73
2:H:67:ILE:HD13	3:I:239:VAL:HA	1.71	0.72
3:C:237:ARG:HH21	3:C:237:ARG:CB	2.02	0.72
3:L:175:GLN:HE22	3:L:177:ARG:HH21	1.38	0.72
3:C:175:GLN:HE22	3:C:177:ARG:HH11	1.37	0.71
2:K:3:SER:O	2:K:7:GLU:HG3	1.90	0.71
2:H:67:ILE:CD1	3:I:239:VAL:HA	2.20	0.71
2:E:148:LEU:HD12	2:E:151:TYR:HB2	1.72	0.71
3:C:237:ARG:NH2	3:C:237:ARG:HB2	2.05	0.70
2:H:147:GLY:C	2:H:148:LEU:HD12	2.12	0.69
2:E:142:VAL:HG22	2:E:148:LEU:HG	1.75	0.69
2:E:124:GLN:HE21	1:G:84:ASN:ND2	1.91	0.68
2:B:148:LEU:CD1	3:C:32:MET:HE1	2.22	0.67
2:H:58:ILE:H	2:H:58:ILE:HD13	1.58	0.67
2:H:142:VAL:N	2:H:148:LEU:HD11	2.10	0.67
3:C:237:ARG:HH21	3:C:237:ARG:HG3	1.59	0.66
1:D:11:ASP:OD2	1:D:14:HIS:HB2	1.96	0.66
2:H:134:LYS:HE3	2:H:138:MET:HE1	1.78	0.66
1:A:86:GLU:H	1:A:86:GLU:CD	1.99	0.65
2:E:24:HIS:ND1	2:K:24:HIS:HD2	1.95	0.64
1:A:111:ASN:O	3:C:175:GLN:HG3	1.97	0.64
2:H:134:LYS:HE3	2:H:138:MET:CE	2.28	0.64
2:K:66:GLU:HB2	2:K:69:GLU:HG3	1.78	0.64
3:F:142:SER:OG	3:F:147:ARG:HD3	1.98	0.63
2:B:66:GLU:HB2	2:B:69:GLU:HG3	1.80	0.63
2:B:124:GLN:HE21	1:J:84:ASN:ND2	1.97	0.62
2:E:67:ILE:HG12	3:F:239:VAL:HB	1.80	0.62
2:E:136:VAL:O	2:E:140:GLU:HG3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:82:ILE:CD1	3:I:117:ALA:HB2	2.28	0.61
2:E:148:LEU:HD21	3:F:32:MET:CE	2.30	0.61
2:H:148:LEU:HA	6:H:233:HOH:O	2.00	0.61
2:E:148:LEU:HD11	3:F:32:MET:CE	2.31	0.60
3:I:161:GLN:HG2	6:L:3108:HOH:O	2.00	0.60
2:B:136:VAL:O	2:B:140:GLU:HG3	2.02	0.60
1:G:111:ASN:O	3:I:175:GLN:HG3	2.02	0.59
3:F:238:PRO:O	3:F:239:VAL:HG23	2.02	0.59
1:G:7:LYS:HB2	1:G:8:PRO:HD3	1.83	0.59
1:A:84:ASN:HD22	1:J:7:LYS:HZ1	1.49	0.58
2:B:148:LEU:HG	3:C:32:MET:CE	2.33	0.58
2:H:138:MET:O	2:H:148:LEU:HD21	2.03	0.58
2:B:142:VAL:HG21	3:C:29:ILE:HD11	1.85	0.58
3:C:164:ALA:HB3	3:F:161:GLN:HE21	1.69	0.57
2:H:66:GLU:HB2	2:H:69:GLU:HG3	1.87	0.57
2:H:136:VAL:O	2:H:140:GLU:HG3	2.05	0.57
3:F:175:GLN:HE22	3:F:177:ARG:HH11	1.52	0.57
2:B:142:VAL:HG22	2:B:148:LEU:HD22	1.87	0.57
2:E:66:GLU:HB2	2:E:69:GLU:HG3	1.86	0.57
1:A:84:ASN:ND2	1:J:7:LYS:NZ	2.53	0.56
1:G:45:ILE:HD12	1:G:45:ILE:N	2.20	0.56
2:B:71:ASN:HD21	3:C:53:LYS:NZ	2.03	0.56
3:C:102:THR:O	3:C:103:GLN:HB2	2.06	0.56
2:B:138:MET:O	2:B:142:VAL:HG23	2.06	0.56
2:H:58:ILE:N	2:H:58:ILE:HD13	2.21	0.56
1:A:26:GLN:HE22	3:C:111:CYS:HB2	1.71	0.55
3:C:32:MET:O	3:C:36:GLU:HG3	2.08	0.54
2:B:42:ALA:O	2:B:45:ARG:HG2	2.08	0.54
3:C:161:GLN:HE21	3:F:165:GLU:HG3	1.73	0.54
2:E:67:ILE:CG1	3:F:239:VAL:HB	2.37	0.54
2:B:152:LEU:HA	6:B:226:HOH:O	2.06	0.53
2:E:148:LEU:CD1	3:F:32:MET:SD	2.96	0.53
3:C:207:GLU:HG2	3:C:225:THR:HG22	1.91	0.53
3:I:224:ILE:N	3:I:224:ILE:HD12	2.23	0.53
1:A:117:GLU:HG3	6:C:3011:HOH:O	2.07	0.53
3:C:127:VAL:HG22	3:C:128:CYS:N	2.24	0.53
3:I:207:GLU:HG2	3:I:225:THR:CG2	2.39	0.53
3:L:127:VAL:HG22	3:L:128:CYS:N	2.24	0.53
1:D:94:ILE:HD12	1:D:94:ILE:N	2.24	0.53
2:K:142:VAL:HG21	3:L:29:ILE:HD11	1.92	0.52
2:E:142:VAL:HG21	3:F:24:VAL:HG21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:127:VAL:HG22	3:F:128:CYS:N	2.24	0.51
3:I:127:VAL:HG22	3:I:128:CYS:N	2.24	0.51
3:C:131:CSD:O	3:C:152:ARG:HD2	2.10	0.51
1:D:79:ASP:HB3	1:D:84:ASN:O	2.10	0.51
2:H:67:ILE:HD11	3:I:239:VAL:HA	1.93	0.51
3:C:75:PRO:O	3:C:79:LYS:HG3	2.10	0.51
1:G:79:ASP:HB3	1:G:84:ASN:O	2.11	0.51
3:I:23:GLU:HG2	3:I:23:GLU:O	2.10	0.51
3:F:102:THR:O	3:F:103:GLN:HB2	2.10	0.51
3:L:131:CSD:O	3:L:152:ARG:HD2	2.10	0.51
6:K:247:HOH:O	3:L:234:LYS:HE3	2.11	0.51
3:I:201:THR:OG1	3:I:204:GLN:HG3	2.11	0.50
2:B:148:LEU:CG	3:C:32:MET:HE1	2.40	0.50
2:B:148:LEU:CG	3:C:32:MET:CE	2.90	0.50
3:C:62:LEU:HB2	3:C:63:PRO:HD3	1.94	0.50
1:A:71:VAL:HG11	1:A:117:GLU:OE2	2.11	0.50
2:E:124:GLN:HE21	1:G:84:ASN:HD21	1.57	0.49
2:B:145:GLY:HA2	6:B:281:HOH:O	2.12	0.49
3:I:62:LEU:HB2	6:I:3045:HOH:O	2.11	0.49
2:H:46:VAL:HG12	2:H:96:CYS:SG	2.53	0.49
3:L:102:THR:O	3:L:103:GLN:HB2	2.13	0.49
1:A:125:LYS:HG3	1:A:125:LYS:O	2.11	0.48
3:L:62:LEU:O	3:L:66:ARG:HG2	2.13	0.48
2:K:116:ALA:O	2:K:119:ILE:HG22	2.13	0.48
3:F:201:THR:OG1	3:F:204:GLN:HG3	2.13	0.48
2:E:28:HIS:CE1	3:I:153:ARG:HH12	2.31	0.48
1:J:12:ARG:HD3	2:K:125:PHE:CZ	2.48	0.48
3:C:207:GLU:HG2	3:C:225:THR:CG2	2.44	0.48
3:I:131:CSD:O	3:I:152:ARG:HD2	2.13	0.48
2:E:45:ARG:NH1	2:E:100:GLN:HE22	2.11	0.48
2:K:5:ILE:N	2:K:5:ILE:HD12	2.29	0.48
2:E:148:LEU:CD1	2:E:151:TYR:HB2	2.42	0.48
2:E:148:LEU:HD13	3:F:32:MET:SD	2.54	0.48
1:G:45:ILE:CD1	1:G:45:ILE:N	2.77	0.47
2:K:46:VAL:CG1	2:K:96:CYS:SG	3.02	0.47
2:B:126:VAL:HG12	2:B:127:THR:O	2.13	0.47
2:B:147:GLY:N	2:B:154:PRO:HA	2.28	0.47
3:F:171:PRO:HB2	3:F:174:VAL:HG23	1.97	0.47
2:B:152:LEU:HD12	2:B:153:PRO:HD2	1.95	0.47
2:H:46:VAL:CG1	2:H:96:CYS:SG	3.02	0.47
3:F:173:GLU:CD	3:F:173:GLU:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:TYR:HA	2:B:46:VAL:HG11	1.96	0.47
3:I:54:ASP:O	3:I:58:THR:HG23	2.15	0.47
3:C:123:LYS:HD3	3:C:123:LYS:C	2.35	0.47
3:F:99:SER:HB2	3:F:100:PRO:HD3	1.97	0.47
2:K:118:ARG:HG3	2:K:122:ASP:OD2	2.15	0.47
3:C:175:GLN:HE22	3:C:177:ARG:NH1	2.09	0.46
3:C:170:LEU:HD13	3:C:176:ILE:HD11	1.97	0.46
2:B:146:GLN:O	2:B:146:GLN:HG2	2.16	0.46
3:C:127:VAL:HB	3:C:155:LEU:HD23	1.97	0.46
2:B:142:VAL:HG22	2:B:148:LEU:HD13	1.97	0.46
2:H:58:ILE:HD12	3:I:150:ASN:CG	2.36	0.46
3:F:126:VAL:HG12	3:F:127:VAL:N	2.30	0.46
3:C:126:VAL:HG12	3:C:127:VAL:N	2.31	0.46
3:C:237:ARG:NH2	3:C:237:ARG:HG3	2.28	0.46
1:D:121:ARG:HG3	1:D:122:TYR:CD1	2.51	0.46
6:B:191:HOH:O	1:J:79:ASP:HA	2.16	0.46
2:B:134:LYS:HE2	3:C:36:GLU:HB2	1.98	0.46
2:E:148:LEU:CD2	3:F:32:MET:SD	3.04	0.46
2:H:67:ILE:HG21	2:H:150:GLU:HG3	1.97	0.46
6:E:250:HOH:O	2:K:24:HIS:HE1	1.98	0.46
2:E:118:ARG:HG3	2:E:122:ASP:OD2	2.17	0.45
1:G:71:VAL:HG11	1:G:117:GLU:OE2	2.16	0.45
3:I:226:VAL:HG23	6:I:3114:HOH:O	2.15	0.45
2:K:5:ILE:HD13	6:K:239:HOH:O	2.14	0.45
1:A:12:ARG:HD3	2:B:125:PHE:CZ	2.52	0.45
1:A:9:VAL:HG23	1:A:9:VAL:O	2.16	0.45
3:F:131:CSD:O	3:F:152:ARG:HD3	2.16	0.45
2:E:144:SER:OG	2:E:146:GLN:HG3	2.16	0.45
2:H:118:ARG:HG3	2:H:122:ASP:OD2	2.17	0.45
1:G:20:THR:HA	3:I:104:PHE:HB3	1.98	0.45
3:I:165:GLU:HG3	3:L:161:GLN:HE21	1.82	0.45
2:K:137:GLU:OE2	2:K:141:ARG:NH2	2.33	0.45
3:C:123:LYS:HG3	3:C:170:LEU:HD11	1.98	0.45
2:E:66:GLU:CD	3:F:235:ALA:HB1	2.38	0.45
1:A:28:PHE:HA	1:A:31:MET:HE3	1.99	0.44
2:E:4:SER:HB3	2:E:7:GLU:OE1	2.17	0.44
3:F:40:GLU:HG3	6:F:3098:HOH:O	2.17	0.44
3:I:106:THR:OG1	3:I:109:ASP:HB2	2.16	0.44
3:L:54:ASP:O	3:L:58:THR:HG23	2.18	0.44
2:H:58:ILE:CD1	2:H:58:ILE:N	2.79	0.44
3:L:66:ARG:NH2	3:L:203:ASP:OD1	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:5:ILE:HD13	6:E:290:HOH:O	2.17	0.44
3:I:118:ASP:OD1	3:I:124:HIS:HD2	2.00	0.44
3:F:47:GLU:O	3:F:51:VAL:HG23	2.18	0.44
2:H:116:ALA:O	2:H:119:ILE:HG22	2.16	0.44
2:K:45:ARG:NH1	2:K:100:GLN:HE22	2.16	0.44
1:A:79:ASP:HA	6:A:3013:HOH:O	2.18	0.44
2:B:80:ALA:HA	2:B:85:TRP:O	2.18	0.44
6:I:3120:HOH:O	3:L:161:GLN:HG2	2.17	0.44
1:D:92:TYR:HB2	1:D:94:ILE:HD11	1.99	0.44
2:H:138:MET:HB3	3:I:29:ILE:HG23	2.00	0.44
1:A:79:ASP:HB3	1:A:84:ASN:O	2.18	0.44
3:F:150:ASN:OD1	3:F:154:ARG:HD3	2.18	0.44
1:A:104:GLU:HG3	6:C:3168:HOH:O	2.18	0.44
2:H:134:LYS:HE3	2:H:138:MET:HE3	2.00	0.44
3:I:207:GLU:HG2	3:I:225:THR:HG22	1.99	0.44
2:B:119:ILE:HD11	2:B:123:LYS:HE3	1.98	0.43
2:E:80:ALA:HA	2:E:85:TRP:O	2.17	0.43
3:F:67:LEU:C	3:F:67:LEU:HD23	2.39	0.43
2:K:120:LEU:HD22	3:L:43:LEU:CD1	2.48	0.43
3:C:164:ALA:HB3	3:F:161:GLN:NE2	2.31	0.43
2:B:142:VAL:HG21	3:C:24:VAL:HG21	2.00	0.43
2:E:116:ALA:O	2:E:119:ILE:HG22	2.19	0.43
3:C:136:ARG:HB2	3:C:137:PRO:HD3	2.01	0.43
2:H:80:ALA:HA	2:H:85:TRP:O	2.19	0.43
3:F:62:LEU:HB3	3:F:63:PRO:HD3	2.00	0.43
3:C:171:PRO:HB2	3:C:174:VAL:HG23	2.01	0.43
2:E:148:LEU:HB3	2:E:152:LEU:HB3	2.01	0.43
3:I:106:THR:HG1	3:I:109:ASP:HB2	1.84	0.43
1:A:119:PRO:HG2	1:A:122:TYR:HD2	1.84	0.42
1:A:12:ARG:NH2	6:A:3019:HOH:O	2.51	0.42
1:G:119:PRO:HG2	1:G:122:TYR:HD2	1.84	0.42
2:H:138:MET:CE	3:I:33:ALA:HA	2.50	0.42
2:H:41:ARG:HG2	2:K:12:HIS:HD2	1.84	0.42
3:F:116:LEU:HB3	3:F:124:HIS:CD2	2.55	0.42
2:H:142:VAL:HG22	2:H:148:LEU:HD13	2.00	0.42
3:L:126:VAL:HG12	3:L:127:VAL:N	2.35	0.42
3:C:54:ASP:O	3:C:58:THR:HG23	2.19	0.42
1:D:84:ASN:ND2	2:H:124:GLN:HE21	2.17	0.42
2:E:46:VAL:CG1	2:E:96:CYS:SG	3.08	0.42
3:F:100:PRO:HA	3:F:101:PRO:HD3	1.91	0.42
3:L:157:ARG:HD3	3:L:158:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:99:SER:HB2	3:C:100:PRO:HD3	2.02	0.42
3:I:132:SER:O	3:I:133:CSO:C	2.68	0.42
2:H:148:LEU:HD12	2:H:148:LEU:N	2.34	0.42
2:E:45:ARG:HH11	2:E:100:GLN:HE22	1.67	0.41
2:B:143:ALA:C	2:B:145:GLY:H	2.24	0.41
2:E:4:SER:HB3	2:E:7:GLU:CG	2.50	0.41
2:H:26:GLN:NE2	6:H:220:HOH:O	2.53	0.41
3:I:103:GLN:HG2	3:I:103:GLN:O	2.20	0.41
3:I:126:VAL:HG12	3:I:127:VAL:N	2.35	0.41
2:B:118:ARG:HG3	2:B:122:ASP:OD2	2.20	0.41
3:C:161:GLN:HG2	6:F:3070:HOH:O	2.21	0.41
1:A:84:ASN:HD22	1:J:7:LYS:CE	2.33	0.41
3:I:67:LEU:C	3:I:67:LEU:HD23	2.41	0.41
3:C:99:SER:N	3:C:100:PRO:CD	2.84	0.41
3:F:125:VAL:HA	3:F:216:GLY:HA3	2.02	0.41
3:C:132:SER:O	3:C:133:CSO:C	2.69	0.41
2:E:148:LEU:N	2:E:152:LEU:O	2.54	0.41
3:L:132:SER:O	3:L:133:CSO:C	2.68	0.41
1:D:50:ASP:O	1:D:51:LEU:C	2.60	0.41
3:C:100:PRO:HA	3:C:101:PRO:HD3	1.96	0.41
3:F:123:LYS:HG3	3:F:170:LEU:HD11	2.03	0.41
2:E:4:SER:HB3	2:E:7:GLU:HG3	2.03	0.40
2:H:148:LEU:HD23	3:I:32:MET:SD	2.61	0.40
3:L:67:LEU:HD23	3:L:67:LEU:C	2.41	0.40
1:D:71:VAL:HG11	1:D:117:GLU:OE2	2.21	0.40
2:E:61:HIS:HE1	6:F:3056:HOH:O	2.04	0.40
3:F:157:ARG:HD3	3:F:158:TRP:CH2	2.56	0.40
3:F:154:ARG:NH2	3:F:165:GLU:OE2	2.43	0.40
2:K:152:LEU:HD12	2:K:153:PRO:HD2	2.03	0.40
1:D:93:SER:C	1:D:94:ILE:HD12	2.42	0.40
2:B:126:VAL:HG13	2:B:130:GLU:HB2	2.03	0.40
3:F:160:ARG:NE	3:F:176:ILE:HD12	2.36	0.40
1:J:111:ASN:O	3:L:175:GLN:HG3	2.21	0.40
1:J:7:LYS:HA	1:J:8:PRO:HD3	1.92	0.40
2:K:127:THR:OG1	2:K:130:GLU:HG3	2.22	0.40
1:A:84:ASN:ND2	1:J:7:LYS:HZ3	2.19	0.40
2:E:127:THR:OG1	2:E:130:GLU:HG3	2.22	0.40
1:J:125:LYS:HG3	6:J:3134:HOH:O	2.21	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%)	111 (95%)	6 (5%)	0	100	100
1	D	117/126 (93%)	108 (92%)	9 (8%)	0	100	100
1	G	118/126 (94%)	114 (97%)	4 (3%)	0	100	100
1	J	118/126 (94%)	114 (97%)	4 (3%)	0	100	100
2	B	150/157 (96%)	144 (96%)	5 (3%)	1 (1%)	24	17
2	E	149/157 (95%)	142 (95%)	7 (5%)	0	100	100
2	H	154/157 (98%)	147 (96%)	7 (4%)	0	100	100
2	K	150/157 (96%)	146 (97%)	4 (3%)	0	100	100
3	C	213/243 (88%)	205 (96%)	8 (4%)	0	100	100
3	F	212/243 (87%)	204 (96%)	8 (4%)	0	100	100
3	I	213/243 (88%)	205 (96%)	8 (4%)	0	100	100
3	L	212/243 (87%)	205 (97%)	7 (3%)	0	100	100
All	All	1923/2104 (91%)	1845 (96%)	77 (4%)	1 (0%)	53	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	144	SER

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	101/108 (94%)	100 (99%)	1 (1%)	78	83
1	D	101/108 (94%)	101 (100%)	0	100	100
1	G	102/108 (94%)	101 (99%)	1 (1%)	78	83
1	J	102/108 (94%)	101 (99%)	1 (1%)	78	83
2	B	130/134 (97%)	128 (98%)	2 (2%)	67	72
2	E	129/134 (96%)	126 (98%)	3 (2%)	53	56
2	H	133/134 (99%)	128 (96%)	5 (4%)	36	33
2	K	130/134 (97%)	128 (98%)	2 (2%)	67	72
3	C	188/212 (89%)	186 (99%)	2 (1%)	76	80
3	F	187/212 (88%)	185 (99%)	2 (1%)	76	80
3	I	188/212 (89%)	187 (100%)	1 (0%)	90	93
3	L	187/212 (88%)	187 (100%)	0	100	100
All	All	1678/1816 (92%)	1658 (99%)	20 (1%)	74	78

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
2	B	48	HIS
2	B	85	TRP
3	C	198	ASP
3	C	237	ARG
2	E	48	HIS
2	E	70	LEU
2	E	85	TRP
3	F	37	LEU
3	F	239	VAL
1	G	125	LYS
2	H	39	LEU
2	H	48	HIS
2	H	58	ILE
2	H	85	TRP
2	H	137	GLU
3	I	62	LEU
1	J	99	LYS
2	K	48	HIS
2	K	85	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	84	ASN
1	A	111	ASN
2	B	26	GLN
2	B	37	HIS
2	B	71	ASN
2	B	100	GLN
3	C	175	GLN
2	E	26	GLN
2	E	61	HIS
2	E	94	GLN
2	E	100	GLN
2	E	146	GLN
3	F	49	HIS
3	F	161	GLN
3	F	175	GLN
1	G	14	HIS
1	G	84	ASN
1	G	87	ASN
2	H	28	HIS
2	H	61	HIS
2	H	100	GLN
2	H	124	GLN
3	I	103	GLN
3	I	112	ASN
3	I	124	HIS
3	I	141	GLN
3	I	169	GLN
3	I	175	GLN
1	J	84	ASN
1	J	87	ASN
2	K	12	HIS
2	K	24	HIS
2	K	26	GLN
2	K	100	GLN
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	4,7,8	1.43	1 (25%)	2,8,10	1.76	1 (50%)
3	CSO	C	133	3,5	4,6,7	1.60	1 (25%)	1,6,8	1.41	0
3	CSD	F	131	3,5	4,7,8	1.58	1 (25%)	2,8,10	1.67	0
3	CSO	F	133	3,5	4,6,7	1.41	1 (25%)	1,6,8	1.47	0
3	CSD	I	131	3,5	4,7,8	1.28	1 (25%)	2,8,10	1.75	1 (50%)
3	CSO	I	133	3,5	4,6,7	1.48	1 (25%)	1,6,8	1.53	0
3	CSD	L	131	3,5	4,7,8	1.38	1 (25%)	2,8,10	1.73	1 (50%)
3	CSO	L	133	3,5	4,6,7	1.56	1 (25%)	1,6,8	1.50	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CSD	C	131	3,5	-	1/2/6/8	0/0/0/0
3	CSO	C	133	3,5	-	0/1/5/7	0/0/0/0
3	CSD	F	131	3,5	-	1/2/6/8	0/0/0/0
3	CSO	F	133	3,5	-	0/1/5/7	0/0/0/0
3	CSD	I	131	3,5	-	1/2/6/8	0/0/0/0
3	CSO	I	133	3,5	-	0/1/5/7	0/0/0/0
3	CSD	L	131	3,5	-	1/2/6/8	0/0/0/0
3	CSO	L	133	3,5	-	0/1/5/7	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	131	CSD	CA-C	2.18	1.53	1.50
3	L	131	CSD	CA-C	2.39	1.53	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	131	CSD	CA-C	2.50	1.53	1.50
3	F	133	CSO	CA-C	2.59	1.53	1.50
3	I	133	CSO	CA-C	2.70	1.53	1.50
3	F	131	CSD	CA-C	2.75	1.53	1.50
3	L	133	CSO	CA-C	2.84	1.54	1.50
3	C	133	CSO	CA-C	2.94	1.54	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	131	CSD	OD1-SG-CB	2.06	109.46	105.54
3	L	131	CSD	OD1-SG-CB	2.08	109.50	105.54
3	C	131	CSD	OD1-SG-CB	2.08	109.50	105.54

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	I	131	CSD	CA-CB-SG-OD1
3	C	131	CSD	CA-CB-SG-OD1
3	L	131	CSD	CA-CB-SG-OD1
3	F	131	CSD	CA-CB-SG-OD1

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	CSO	1	0
3	F	131	CSD	1	0
3	I	131	CSD	1	0
3	I	133	CSO	1	0
3	L	131	CSD	1	0
3	L	133	CSO	1	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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	SO4	A	3005	-	4,4,4	0.36	0	6,6,6	0.10	0
4	SO4	C	3001	-	4,4,4	0.30	0	6,6,6	0.10	0
4	SO4	D	3006	-	4,4,4	0.34	0	6,6,6	0.08	0
4	SO4	F	3002	-	4,4,4	0.32	0	6,6,6	0.10	0
4	SO4	G	3007	-	4,4,4	0.34	0	6,6,6	0.08	0
4	SO4	I	3003	-	4,4,4	0.29	0	6,6,6	0.09	0
4	SO4	I	3011	-	4,4,4	0.38	0	6,6,6	0.09	0
4	SO4	I	3012	-	4,4,4	0.44	0	6,6,6	0.14	0
4	SO4	J	3008	-	4,4,4	0.37	0	6,6,6	0.07	0
4	SO4	L	3004	-	4,4,4	0.34	0	6,6,6	0.06	0
4	SO4	L	3009	-	4,4,4	0.33	0	6,6,6	0.09	0
4	SO4	L	3010	-	4,4,4	0.38	0	6,6,6	0.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SO4	A	3005	-	-	0/0/0/0	0/0/0/0
4	SO4	C	3001	-	-	0/0/0/0	0/0/0/0
4	SO4	D	3006	-	-	0/0/0/0	0/0/0/0
4	SO4	F	3002	-	-	0/0/0/0	0/0/0/0
4	SO4	G	3007	-	-	0/0/0/0	0/0/0/0
4	SO4	I	3003	-	-	0/0/0/0	0/0/0/0
4	SO4	I	3011	-	-	0/0/0/0	0/0/0/0
4	SO4	I	3012	-	-	0/0/0/0	0/0/0/0
4	SO4	J	3008	-	-	0/0/0/0	0/0/0/0
4	SO4	L	3004	-	-	0/0/0/0	0/0/0/0
4	SO4	L	3009	-	-	0/0/0/0	0/0/0/0
4	SO4	L	3010	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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%)	-0.31	2 (1%) 70 68	12, 19, 32, 40	0
1	D	119/126 (94%)	0.10	7 (5%) 22 22	17, 26, 42, 50	0
1	G	120/126 (95%)	-0.02	6 (5%) 29 28	13, 21, 37, 57	0
1	J	120/126 (95%)	-0.22	2 (1%) 70 68	11, 19, 34, 43	0
2	B	152/157 (96%)	0.03	11 (7%) 15 15	11, 18, 46, 57	0
2	E	151/157 (96%)	0.01	9 (5%) 22 21	13, 20, 49, 58	0
2	H	156/157 (99%)	0.35	14 (8%) 9 9	12, 20, 55, 67	0
2	K	152/157 (96%)	-0.44	1 (0%) 87 87	10, 16, 29, 43	0
3	C	215/243 (88%)	-0.03	7 (3%) 46 46	10, 22, 37, 49	0
3	F	214/243 (88%)	0.09	5 (2%) 60 59	15, 24, 35, 48	0
3	I	215/243 (88%)	-0.16	4 (1%) 66 65	10, 17, 30, 47	0
3	L	214/243 (88%)	-0.34	1 (0%) 90 90	9, 16, 26, 33	0
All	All	1947/2104 (92%)	-0.08	69 (3%) 44 43	9, 20, 36, 67	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	156	ALA	12.4
2	H	157	LYS	10.8
3	F	239	VAL	8.2
2	H	3	SER	8.1
2	H	148	LEU	8.0
2	H	154	PRO	7.2
2	H	2	SER	7.2
2	B	153	PRO	7.1
2	E	153	PRO	7.1
2	B	152	LEU	7.1
2	B	154	PRO	7.0

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Mol	Chain	Res	Type	RSRZ
2	B	148	LEU	6.9
2	B	144	SER	6.5
2	H	152	LEU	6.1
2	E	154	PRO	6.0
2	H	153	PRO	5.8
2	E	148	LEU	5.8
2	H	147	GLY	5.4
2	H	144	SER	5.3
2	E	147	GLY	5.2
2	H	155	LYS	5.2
2	K	3	SER	4.9
2	B	147	GLY	4.9
2	E	144	SER	4.9
1	A	126	ALA	4.8
3	C	198	ASP	4.6
1	G	126	ALA	4.5
2	B	146	GLN	4.3
3	C	23	GLU	4.1
2	E	145	GLY	4.1
2	E	152	LEU	4.1
1	G	7	LYS	4.1
3	F	76	GLU	3.6
2	B	3	SER	3.6
2	E	146	GLN	3.6
2	H	146	GLN	3.5
1	D	9	VAL	3.5
1	D	126	ALA	3.3
2	H	145	GLY	3.1
3	F	198	ASP	3.0
1	D	8	PRO	2.9
1	A	125	LYS	2.9
1	D	125	LYS	2.9
3	C	173	GLU	2.8
2	B	143	ALA	2.8
2	H	4	SER	2.8
2	B	145	GLY	2.7
1	J	34	LYS	2.6
3	C	199	GLY	2.6
1	G	125	LYS	2.5
3	I	23	GLU	2.5
1	G	9	VAL	2.4
1	D	17	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	G	8	PRO	2.3
3	I	199	GLY	2.3
3	C	234	LYS	2.2
3	F	79	LYS	2.2
1	J	126	ALA	2.2
3	F	127	VAL	2.2
2	E	143	ALA	2.2
2	B	4	SER	2.1
1	G	34	LYS	2.1
3	L	66	ARG	2.1
3	C	200	TRP	2.1
1	D	11	ASP	2.1
1	D	86	GLU	2.1
3	I	76	GLU	2.1
3	I	130	LEU	2.0
3	C	229	LYS	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CSO	F	133	7/8	0.80	0.21	13,14,18,19	1
3	CSO	I	133	7/8	0.85	0.21	10,12,17,18	1
3	CSD	F	131	8/9	0.88	0.21	11,14,15,17	2
3	CSD	C	131	8/9	0.93	0.19	8,11,13,15	2
3	CSD	I	131	8/9	0.94	0.21	8,10,13,14	2
3	CSO	C	133	7/8	0.94	0.13	9,11,16,16	1
3	CSO	L	133	7/8	0.96	0.16	6,9,13,15	1
3	CSD	L	131	8/9	0.97	0.17	5,9,10,12	2

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

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(\AA^2)	Q<0.9
5	3CO	F	301	1/1	-	-	18,18,18,18	1
4	SO4	F	3002	5/5	0.97	0.14	37,37,40,40	0
5	3CO	I	301	1/1	-	-	13,13,13,13	1
5	3CO	C	301	1/1	-	-	14,14,14,14	1
4	SO4	L	3010	5/5	0.92	0.21	49,49,50,52	0
4	SO4	I	3011	5/5	0.94	0.19	49,49,50,52	0
4	SO4	A	3005	5/5	0.95	0.12	42,42,44,45	0
4	SO4	L	3009	5/5	0.96	0.14	47,48,49,50	0
4	SO4	I	3003	5/5	0.97	0.16	33,34,35,36	0
4	SO4	I	3012	5/5	0.97	0.12	51,51,51,53	0
4	SO4	J	3008	5/5	0.98	0.11	46,46,47,47	0
4	SO4	D	3006	5/5	0.98	0.09	37,37,38,39	0
4	SO4	C	3001	5/5	0.98	0.10	32,34,36,37	0
5	3CO	L	301	1/1	-	-	12,12,12,12	1
4	SO4	G	3007	5/5	0.97	0.09	41,41,41,42	0
4	SO4	L	3004	5/5	0.97	0.14	30,34,37,37	0

6.5 Other polymers

There are no such residues in this entry.