



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 11:49 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 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/ValidationPDFNotes.html>

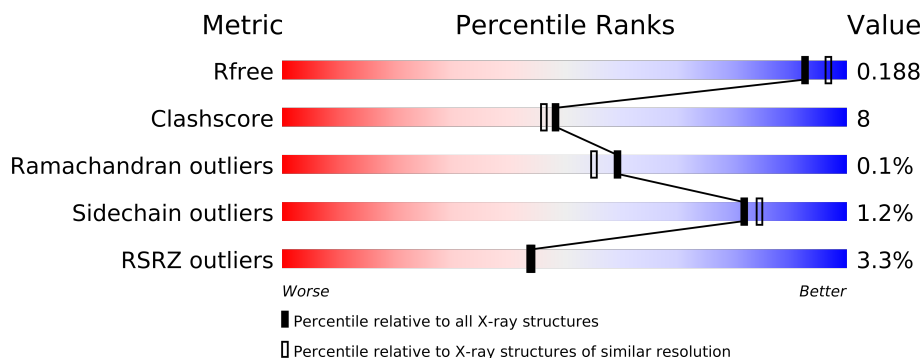
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	126	
1	D	126	
1	G	126	
1	J	126	
2	B	157	
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 that are outliers for geometric or electron-

density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	SO4	I	3011	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17962 atoms, of which 0 are hydrogen and 0 are deuterium.

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

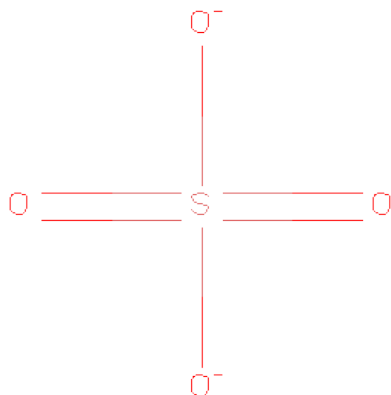
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	1	Total	Co	1	0
			1	1		
4	L	1	Total	Co	1	0
			1	1		
4	C	1	Total	Co	1	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	1	Total	Co	1	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	I	1	Total	O	S	0	0
			5	4	1		
5	I	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

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

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: Thiocyanate hydrolase alpha subunit

Chain A: 



- Molecule 1: Thiocyanate hydrolase alpha subunit

Chain D: 



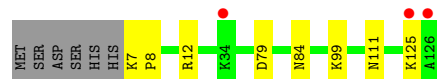
- Molecule 1: Thiocyanate hydrolase alpha subunit

Chain G: 



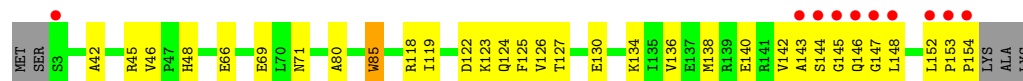
- Molecule 1: Thiocyanate hydrolase alpha subunit

Chain J: 



- Molecule 2: Thiocyanate hydrolase beta subunit

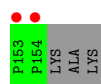
Chain B: 



- Molecule 2: Thiocyanate hydrolase beta subunit

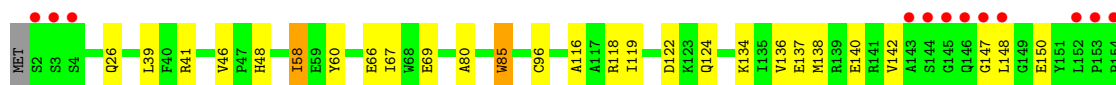
Chain E: 





- Molecule 2: Thiocyanate hydrolase beta subunit

Chain H:



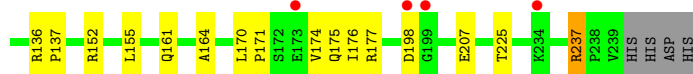
- Molecule 2: Thiocyanate hydrolase beta subunit

Chain K:



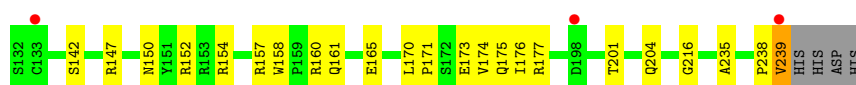
- Molecule 3: Thiocyanate hydrolase gamma subunit

Chain C:



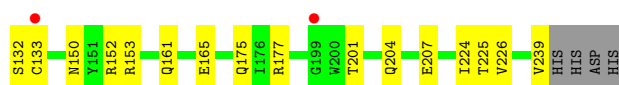
- Molecule 3: Thiocyanate hydrolase gamma subunit

Chain F:



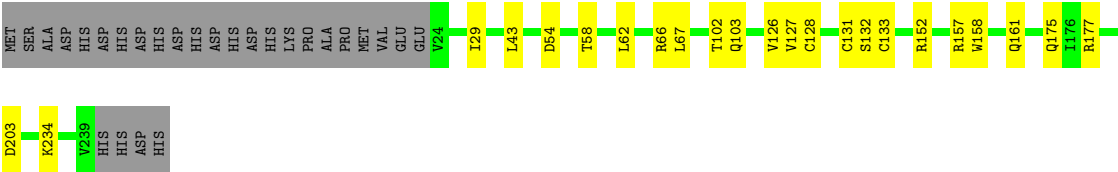
- Molecule 3: Thiocyanate hydrolase gamma subunit

Chain I:



- Molecule 3: Thiocyanate hydrolase gamma subunit

Chain L: 



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 0.187 , 0.188	Depositor DCC
R_{free} test set	14043 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 43.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 278076 reflections	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.

5 Model quality

5.1 Standard geometry

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	965	0	914	23	0
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	C	1	0	0	0	0
4	F	1	0	0	0	0
4	I	1	0	0	0	0
4	L	1	0	0	0	0
5	A	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
5	F	5	0	0	0	0
5	G	5	0	0	0	0
5	I	15	0	0	0	0
5	J	5	0	0	0	0
5	L	15	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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (244) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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
2:H:134:LYS:HE3	2:H:138:MET:HE1	1.78	0.66
1:D:11:ASP:OD2	1:D:14:HIS:HB2	1.96	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
2:H:134:LYS:HE3	2:H:138:MET:CE	2.28	0.64
1:A:111:ASN:O	3:C:175:GLN:HG3	1.97	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
3:I:82:ILE:CD1	3:I:117:ALA:HB2	2.28	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:H:138:MET:O	2:H:148:LEU:HD21	2.03	0.58
2:B:148:LEU:HG	3:C:32:MET:CE	2.33	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
2:E:67:ILE:CG1	3:F:239:VAL:HB	2.37	0.54
3:C:161:GLN:HE21	3:F:165:GLU:HG3	1.73	0.54
2:E:148:LEU:CD1	3:F:32:MET:SD	2.96	0.53
2:B:152:LEU:HA	6:B:226:HOH:O	2.06	0.53
3:I:224:ILE:N	3:I:224:ILE:HD12	2.23	0.53
3:C:207:GLU:HG2	3:C:225:THR:HG22	1.91	0.53
1:A:117:GLU:HG3	6:C:3011:HOH:O	2.07	0.53
3:L:127:VAL:HG22	3:L:128:CYS:N	2.24	0.53
3:I:207:GLU:HG2	3:I:225:THR:CG2	2.39	0.53
3:C:127:VAL:HG22	3:C: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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:C:131:CSD:O	3:C:152:ARG:HD2	2.10	0.51
2:H:67:ILE:HD11	3:I:239:VAL:HA	1.93	0.51
1:D:79:ASP:HB3	1:D:84:ASN:O	2.10	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: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:F:102:THR:O	3:F:103:GLN:HB2	2.10	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
3:I:62:LEU:HB2	6:I:3045:HOH:O	2.11	0.49
2:B:145:GLY:HA2	6:B:281:HOH:O	2.12	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
3:L:62:LEU:O	3:L:66:ARG:HG2	2.13	0.48
1:A:125:LYS:O	1:A:125:LYS:HG3	2.11	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:K:5:ILE:N	2:K:5:ILE:HD12	2.29	0.48
2:E:45:ARG:NH1	2:E:100:GLN:HE22	2.11	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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:237:ARG:NH2	3:C:237:ARG:HG3	2.28	0.46
3:C:126:VAL:HG12	3:C:127:VAL:N	2.31	0.46
6:B:191:HOH:O	1:J:79:ASP:HA	2.16	0.46
1:D:121:ARG:HG3	1:D:122:TYR:CD1	2.51	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:K:5:ILE:HD13	6:K:239:HOH:O	2.14	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:E:118:ARG:HG3	2:E:122:ASP:OD2	2.17	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
2:E:66:GLU:CD	3:F:235:ALA:HB1	2.38	0.45
3:C:123:LYS:HG3	3:C:170:LEU:HD11	1.98	0.45
3:I:106:THR:OG1	3:I:109:ASP:HB2	2.16	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
1:A:28:PHE:HA	1:A:31:MET:HE3	1.99	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
3:I:118:ASP:OD1	3:I:124:HIS:HD2	2.00	0.44
2:E:5:ILE:HD13	6:E:290:HOH:O	2.17	0.44
2:H:116:ALA:O	2:H:119:ILE:HG22	2.16	0.44
3:F:47:GLU:O	3:F:51:VAL:HG23	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
6:L:3120:HOH:O	3:L:161:GLN:HG2	2.17	0.44
2:B:80:ALA:HA	2:B:85:TRP:O	2.18	0.44
2:H:138:MET:HB3	3:I:29:ILE:HG23	2.00	0.44
1:D:92:TYR:HB2	1:D:94:ILE:HD11	1.99	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
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
1:A:104:GLU:HG3	6:C:3168:HOH:O	2.18	0.44
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:B:119:ILE:HD11	2:B:123:LYS:HE3	1.98	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
2:H:80:ALA:HA	2:H:85:TRP:O	2.19	0.43
3:C:136:ARG:HB2	3:C:137:PRO:HD3	2.01	0.43
3:F:62:LEU:HB3	3:F:63:PRO:HD3	2.00	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
3:C:171:PRO:HB2	3:C:174:VAL:HG23	2.01	0.43
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
1:A:119:PRO:HG2	1:A: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
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:F:116:LEU:HB3	3:F:124:HIS:CD2	2.55	0.42
1:D:84:ASN:ND2	2:H:124:GLN:HE21	2.17	0.42
3:F:100:PRO:HA	3:F:101:PRO:HD3	1.91	0.42
2:E:46:VAL:CG1	2:E:96:CYS:SG	3.08	0.42
3:L:157:ARG:HD3	3:L:158:TRP:CH2	2.54	0.42
3:C:54:ASP:O	3:C:58:THR:HG23	2.19	0.42
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
3:I:126:VAL:HG12	3:I:127:VAL:N	2.35	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
3:I:103:GLN:HG2	3:I:103:GLN:O	2.20	0.41
2:H:26:GLN:NE2	6:H:220:HOH:O	2.53	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
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
3:C:132:SER:O	3:C:133:CSO:C	2.69	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:H:148:LEU:HD23	3:I:32:MET:SD	2.61	0.40
2:E:4:SER:HB3	2:E:7:GLU:HG3	2.03	0.40
3:L:67:LEU:HD23	3:L:67:LEU:C	2.41	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
3:F:157:ARG:HD3	3:F:158:TRP:CH2	2.56	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
1:D:93:SER:C	1:D:94:ILE:HD12	2.42	0.40
1:J:7:LYS:HA	1:J:8:PRO:HD3	1.92	0.40
1:J:111:ASN:O	3:L:175:GLN:HG3	2.21	0.40
2:B:126:VAL:HG13	2:B:130:GLU:HB2	2.03	0.40
2:K:127:THR:OG1	2:K:130:GLU:HG3	2.22	0.40
3:F:160:ARG:NE	3:F:176:ILE:HD12	2.36	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%)	30	20
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%)	59	55

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%)	85	88
1	D	101/108 (94%)	101 (100%)	0	100	100
1	G	102/108 (94%)	101 (99%)	1 (1%)	85	88
1	J	102/108 (94%)	101 (99%)	1 (1%)	85	88
2	B	130/134 (97%)	128 (98%)	2 (2%)	76	79
2	E	129/134 (96%)	126 (98%)	3 (2%)	63	63

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	133/134 (99%)	128 (96%)	5 (4%)	44	39
2	K	130/134 (97%)	128 (98%)	2 (2%)	76	79
3	C	188/212 (89%)	186 (99%)	2 (1%)	84	86
3	F	187/212 (88%)	185 (99%)	2 (1%)	84	86
3	I	188/212 (89%)	187 (100%)	1 (0%)	94	96
3	L	187/212 (88%)	187 (100%)	0	100	100
All	All	1678/1816 (92%)	1658 (99%)	20 (1%)	82	84

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

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Mol	Chain	Res	Type
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 chains 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	3,4	7,7,8	7.00	3 (42%)	6,8,10	3.63	4 (66%)
3	CSO	C	133	3,4	6,6,7	7.65	3 (50%)	3,6,8	0.85	0
3	CSD	F	131	3,4	7,7,8	7.41	3 (42%)	6,8,10	3.52	4 (66%)
3	CSO	F	133	3,4	6,6,7	7.39	3 (50%)	3,6,8	1.04	0
3	CSD	I	131	3,4	7,7,8	7.01	3 (42%)	6,8,10	3.11	3 (50%)
3	CSO	I	133	3,4	6,6,7	7.59	3 (50%)	3,6,8	0.95	0
3	CSD	L	131	3,4	7,7,8	7.05	3 (42%)	6,8,10	3.57	3 (50%)
3	CSO	L	133	3,4	6,6,7	7.53	3 (50%)	3,6,8	1.01	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	3,4	-	1/3/6/8	0/0/0/0
3	CSO	C	133	3,4	-	0/2/5/7	0/0/0/0
3	CSD	F	131	3,4	-	1/3/6/8	0/0/0/0
3	CSO	F	133	3,4	-	0/2/5/7	0/0/0/0
3	CSD	I	131	3,4	-	1/3/6/8	0/0/0/0
3	CSO	I	133	3,4	-	0/2/5/7	0/0/0/0
3	CSD	L	131	3,4	-	1/3/6/8	0/0/0/0
3	CSO	L	133	3,4	-	0/2/5/7	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	131	CSD	O-C	19.03	1.24	1.11
3	C	133	CSO	O-C	18.15	1.23	1.11
3	L	131	CSD	O-C	18.06	1.23	1.11
3	I	133	CSO	O-C	18.03	1.23	1.11
3	I	131	CSD	O-C	18.00	1.23	1.11
3	C	131	CSD	O-C	17.93	1.23	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	133	CSO	O-C	17.86	1.23	1.11
3	F	133	CSO	O-C	17.52	1.23	1.11
3	L	131	CSD	OD1-SG	-3.39	1.43	1.47
3	C	131	CSD	OD1-SG	-3.32	1.43	1.47
3	I	131	CSD	OD1-SG	-3.26	1.43	1.47
3	F	131	CSD	OD1-SG	-3.23	1.43	1.47
3	C	133	CSO	CA-C	3.08	1.54	1.48
3	L	133	CSO	CA-C	3.02	1.54	1.48
3	C	133	CSO	OD-SG	-3.01	1.49	1.62
3	F	133	CSO	OD-SG	-3.00	1.49	1.62
3	I	133	CSO	OD-SG	-2.97	1.49	1.62
3	L	133	CSO	OD-SG	-2.96	1.50	1.62
3	F	131	CSD	CA-C	2.96	1.53	1.48
3	I	133	CSO	CA-C	2.93	1.53	1.48
3	F	133	CSO	CA-C	2.85	1.53	1.48
3	C	131	CSD	CA-C	2.79	1.53	1.48
3	L	131	CSD	CA-C	2.72	1.53	1.48
3	I	131	CSD	CA-C	2.58	1.53	1.48

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	131	CSD	C-CA-N	7.42	121.24	113.83
3	C	131	CSD	C-CA-N	7.42	121.24	113.83
3	F	131	CSD	C-CA-N	7.23	121.05	113.83
3	I	131	CSD	C-CA-N	6.16	119.99	113.83
3	L	131	CSD	OD2-SG-CB	3.12	109.97	97.42
3	C	131	CSD	OD2-SG-CB	3.09	109.87	97.42
3	I	131	CSD	OD2-SG-CB	3.08	109.82	97.42
3	F	131	CSD	OD2-SG-CB	3.08	109.81	97.42
3	C	131	CSD	CA-CB-SG	2.71	114.68	110.82
3	F	131	CSD	CA-CB-SG	2.38	114.22	110.82
3	L	131	CSD	CA-CB-SG	2.26	114.04	110.82
3	C	131	CSD	CB-CA-N	2.12	113.56	110.16
3	F	131	CSD	CB-CA-N	2.11	113.55	110.16
3	I	131	CSD	CA-CB-SG	2.08	113.78	110.82

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

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

There are no ring outliers.

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	$\# Z > 2$
5	SO4	A	3005	-	4,4,4	0.29	0	6,6,6	0.08	0
5	SO4	C	3001	-	4,4,4	0.19	0	6,6,6	0.10	0
5	SO4	D	3006	-	4,4,4	0.27	0	6,6,6	0.06	0
5	SO4	F	3002	-	4,4,4	0.19	0	6,6,6	0.09	0
5	SO4	G	3007	-	4,4,4	0.27	0	6,6,6	0.06	0
5	SO4	I	3003	-	4,4,4	0.18	0	6,6,6	0.09	0
5	SO4	I	3011	-	4,4,4	0.34	0	6,6,6	0.08	0
5	SO4	I	3012	-	4,4,4	0.43	0	6,6,6	0.14	0
5	SO4	J	3008	-	4,4,4	0.28	0	6,6,6	0.07	0
5	SO4	L	3004	-	4,4,4	0.22	0	6,6,6	0.06	0
5	SO4	L	3009	-	4,4,4	0.24	0	6,6,6	0.07	0
5	SO4	L	3010	-	4,4,4	0.32	0	6,6,6	0.08	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
5	SO4	A	3005	-	-	0/0/0/0	0/0/0/0
5	SO4	C	3001	-	-	0/0/0/0	0/0/0/0
5	SO4	D	3006	-	-	0/0/0/0	0/0/0/0
5	SO4	F	3002	-	-	0/0/0/0	0/0/0/0
5	SO4	G	3007	-	-	0/0/0/0	0/0/0/0
5	SO4	I	3003	-	-	0/0/0/0	0/0/0/0
5	SO4	I	3011	-	-	0/0/0/0	0/0/0/0
5	SO4	I	3012	-	-	0/0/0/0	0/0/0/0
5	SO4	J	3008	-	-	0/0/0/0	0/0/0/0
5	SO4	L	3004	-	-	0/0/0/0	0/0/0/0
5	SO4	L	3009	-	-	0/0/0/0	0/0/0/0
5	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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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%)	-0.34	2 (1%) 67 67	12, 19, 32, 40	0
1	D	119/126 (94%)	0.06	5 (4%) 35 34	17, 26, 42, 50	0
1	G	120/126 (95%)	-0.08	5 (4%) 35 34	13, 21, 37, 57	0
1	J	120/126 (95%)	-0.28	3 (2%) 54 54	11, 19, 34, 43	0
2	B	152/157 (96%)	-0.06	10 (6%) 18 17	11, 18, 46, 57	0
2	E	151/157 (96%)	-0.06	8 (5%) 25 25	13, 20, 49, 58	0
2	H	156/157 (99%)	0.29	15 (9%) 8 8	12, 20, 55, 67	0
2	K	152/157 (96%)	-0.45	1 (0%) 84 85	10, 16, 29, 43	0
3	C	217/243 (89%)	-0.07	5 (2%) 57 57	10, 22, 37, 49	2 (0%)
3	F	216/243 (88%)	0.08	7 (3%) 45 45	15, 24, 35, 48	2 (0%)
3	I	217/243 (89%)	-0.19	3 (1%) 72 72	10, 17, 30, 47	2 (0%)
3	L	216/243 (88%)	-0.36	0 100 100	9, 16, 26, 33	2 (0%)
All	All	1955/2104 (92%)	-0.12	64 (3%) 44 44	9, 20, 36, 67	8 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	156	ALA	12.3
2	H	157	LYS	10.6
2	H	148	LEU	8.4
3	F	239	VAL	8.0
2	H	3	SER	7.4
2	H	2	SER	6.6
2	H	154	PRO	6.6
2	B	152	LEU	6.1
2	B	148	LEU	6.0
2	B	154	PRO	5.9
2	B	153	PRO	5.9

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Mol	Chain	Res	Type	RSRZ
2	H	152	LEU	5.8
2	B	144	SER	5.8
2	E	153	PRO	5.6
2	H	147	GLY	5.3
2	E	154	PRO	5.3
2	H	153	PRO	5.2
2	H	155	LYS	5.2
2	E	148	LEU	5.1
2	E	147	GLY	4.9
2	H	144	SER	4.6
2	K	3	SER	4.5
2	E	144	SER	4.3
3	C	198	ASP	4.3
1	A	126	ALA	4.2
1	G	126	ALA	4.1
2	B	147	GLY	4.0
3	C	23	GLU	4.0
2	B	3	SER	3.8
2	E	152	LEU	3.6
1	G	7	LYS	3.5
2	E	146	GLN	3.4
1	D	9	VAL	3.4
2	B	146	GLN	3.4
3	F	76	GLU	3.4
2	E	145	GLY	3.4
1	D	126	ALA	3.2
2	H	146	GLN	3.1
1	D	125	LYS	3.0
2	H	145	GLY	3.0
1	D	8	PRO	2.9
1	A	125	LYS	2.6
3	F	131	CSD	2.6
3	F	198	ASP	2.6
3	C	199	GLY	2.6
3	C	173	GLU	2.6
1	J	34	LYS	2.4
1	G	9	VAL	2.4
1	D	17	LYS	2.4
2	B	143	ALA	2.3
1	G	125	LYS	2.3
3	F	79	LYS	2.3
2	B	145	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
2	H	4	SER	2.2
3	I	23	GLU	2.2
3	I	199	GLY	2.2
1	G	8	PRO	2.2
1	J	125	LYS	2.1
3	F	127	VAL	2.1
3	F	133	CSO	2.1
3	I	133	CSO	2.1
2	H	143	ALA	2.1
3	C	234	LYS	2.1
1	J	126	ALA	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	CSD	C	131	8/9	0.18	2.98	8,11,13,15	2
3	CSD	F	131	8/9	0.21	2.45	11,14,15,17	2
3	CSO	I	133	7/8	0.21	2.37	10,12,17,18	1
3	CSO	F	133	7/8	0.21	2.33	13,14,18,19	1
3	CSD	L	131	8/9	0.16	1.81	5,9,10,12	2
3	CSD	I	131	8/9	0.19	1.30	8,10,13,14	2
3	CSO	L	133	7/8	0.14	1.24	6,9,13,15	1
3	CSO	C	133	7/8	0.12	-0.30	9,11,16,16	1

6.3 Carbohydrates ⓘ

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. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	SO4	I	3011	5/5	0.20	2.95	49,49,50,52	0
5	SO4	L	3010	5/5	0.20	1.53	49,49,50,52	0
5	SO4	L	3004	5/5	0.14	1.47	30,34,37,37	0
5	SO4	I	3012	5/5	0.10	0.94	51,51,51,53	0
5	SO4	I	3003	5/5	0.16	0.38	33,34,35,36	0
5	SO4	A	3005	5/5	0.11	0.19	42,42,44,45	0
5	SO4	F	3002	5/5	0.14	-0.14	37,37,40,40	0
5	SO4	L	3009	5/5	0.12	-0.31	47,48,49,50	0
5	SO4	C	3001	5/5	0.10	-0.56	32,34,36,37	0
5	SO4	D	3006	5/5	0.09	-0.74	37,37,38,39	0
5	SO4	J	3008	5/5	0.08	-1.05	46,46,47,47	0
5	SO4	G	3007	5/5	0.08	-1.18	41,41,41,42	0
4	3CO	C	301	1/1	-	-	14,14,14,14	1
4	3CO	I	301	1/1	-	-	13,13,13,13	1
4	3CO	L	301	1/1	-	-	12,12,12,12	1
4	3CO	F	301	1/1	-	-	18,18,18,18	1

6.5 Other polymers ⓘ

There are no such residues in this entry.