



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:01 AM UTC

PDB ID : 1UQR / pdb_00001uqr
Title : Type II 3-dehydroquinate dehydratase (DHQase) from *Actinobacillus pleuropneumoniae*
Authors : Maes, D.; Gonzalez-Ramirez, L.A.; Lopez-Jaramillo, J.; Yu, B.; De Bondt, H.; Zegers, I.; Afonina, E.; Garcia-Ruiz, J.M.; Gulnik, S.
Deposited on : 2003-10-16
Resolution : 1.70 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

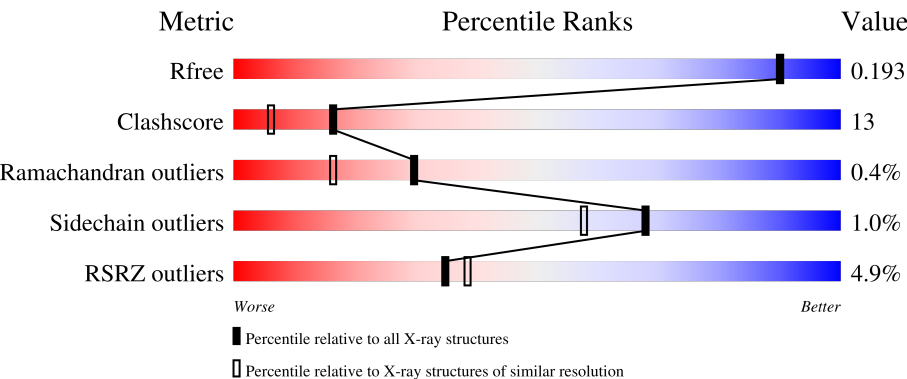
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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

Mol	Chain	Length	Quality of chain
1	A	154	6% 73% 21% • 5%
1	B	154	2% 75% 14% • 10%
1	C	154	3% 73% 23% • •
1	D	154	7% 73% 23% • • •
1	E	154	6% 72% 21% 7%

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Mol	Chain	Length	Quality of chain
1	F	154	5% 77% 18% 5%
1	G	154	3% 68% 23% • 9%
1	H	154	3% 75% 21% •
1	I	154	5% 68% 22% •• 8%
1	J	154	7% 79% 14% • 5%
1	K	154	6% 79% 18% ••
1	L	154	3% 69% 20% • 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16147 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 3-DEHYDROQUINATE DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	1	0
			1156	736	201	214	5			
1	B	138	Total	C	N	O	S	0	1	0
			1089	694	188	202	5			
1	C	148	Total	C	N	O	S	0	1	1
			1161	735	203	216	7			
1	D	152	Total	C	N	O	S	0	1	1
			1195	759	208	222	6			
1	E	143	Total	C	N	O	S	0	1	0
			1130	720	196	209	5			
1	F	146	Total	C	N	O	S	0	1	1
			1145	726	200	213	6			
1	G	140	Total	C	N	O	S	0	1	0
			1106	703	193	205	5			
1	H	148	Total	C	N	O	S	0	1	1
			1161	735	203	216	7			
1	I	141	Total	C	N	O	S	0	1	0
			1115	708	194	208	5			
1	J	146	Total	C	N	O	S	0	1	0
			1156	736	201	214	5			
1	K	152	Total	C	N	O	S	0	1	1
			1195	759	208	222	6			
1	L	138	Total	C	N	O	S	0	1	0
			1089	694	188	202	5			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	K	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		
2	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			8	4	1	3		
3	E	1	Total	C	N	O	0	0
			8	4	1	3		
3	H	1	Total	C	N	O	0	0
			8	4	1	3		
3	K	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	195	Total	O	0	0
			195	195		
4	B	204	Total	O	0	0
			204	204		
4	C	216	Total	O	0	0
			216	216		
4	D	203	Total	O	0	0
			203	203		
4	E	162	Total	O	0	0
			162	162		
4	F	216	Total	O	0	0
			216	216		
4	G	156	Total	O	0	0
			156	156		
4	H	199	Total	O	0	0
			199	199		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	179	Total 179	O 179	0	0
4	J	152	Total 152	O 152	0	0
4	K	199	Total 199	O 199	0	0
4	L	181	Total 181	O 181	0	0

● Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain E: 



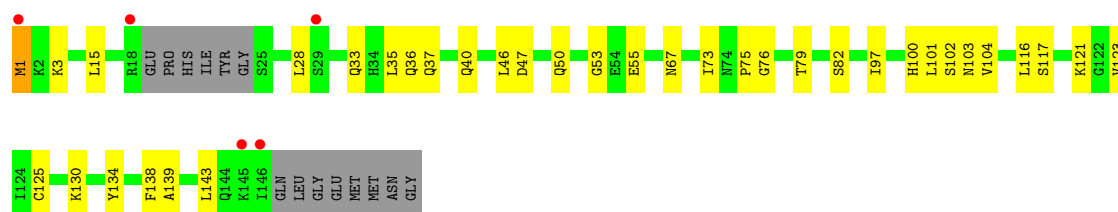
● Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain F: 



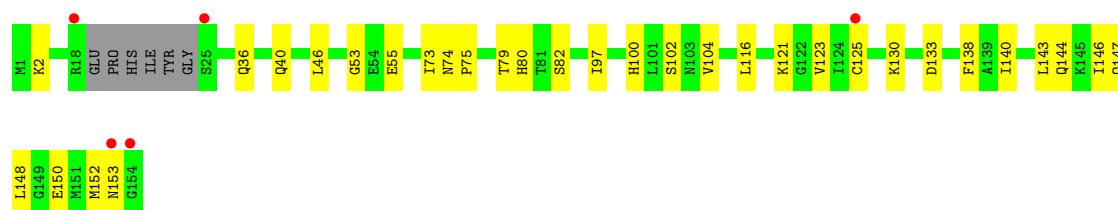
● Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain G: 



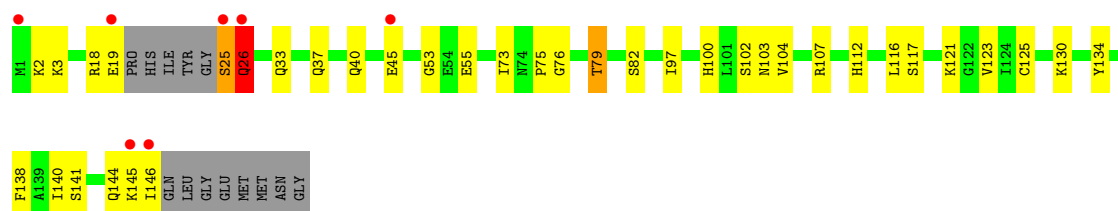
● Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

Chain H: 

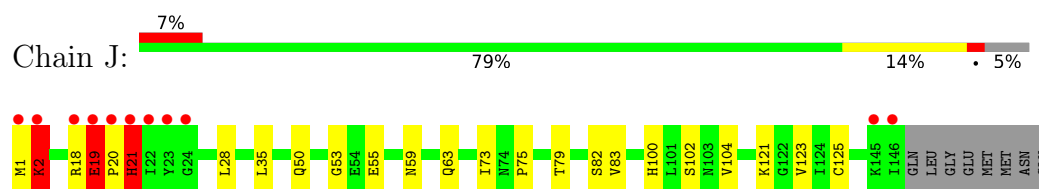


● Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE

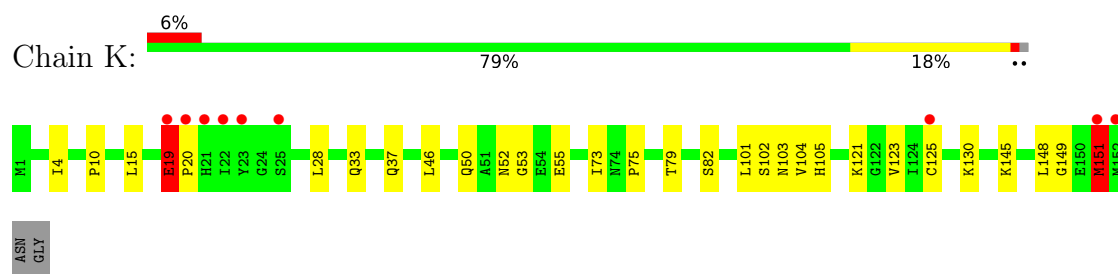
Chain I: 



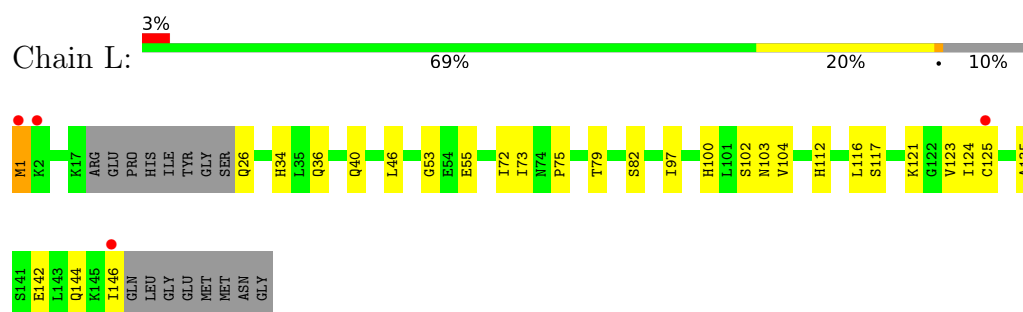
- Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE



- Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE



- Molecule 1: 3-DEHYDROQUINATE DEHYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	129.09Å 131.33Å 161.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.70 20.00 – 1.71	Depositor EDS
% Data completeness (in resolution range)	98.0 (20.00-1.70) 98.3 (20.00-1.71)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.88 (at 1.71Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.178 , 0.199 0.173 , 0.193	Depositor DCC
R_{free} test set	14401 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16147	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, 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	0.41	0/1189	0.90	4/1609 (0.2%)
1	B	0.43	0/1118	0.88	3/1511 (0.2%)
1	C	0.42	1/1190 (0.1%)	0.86	5/1606 (0.3%)
1	D	0.42	1/1228 (0.1%)	0.85	3/1661 (0.2%)
1	E	0.37	0/1160	0.85	3/1567 (0.2%)
1	F	0.41	1/1174 (0.1%)	0.88	4/1585 (0.3%)
1	G	0.40	0/1135	0.84	4/1533 (0.3%)
1	H	0.42	1/1190 (0.1%)	0.85	4/1606 (0.2%)
1	I	0.46	1/1144 (0.1%)	0.91	4/1545 (0.3%)
1	J	0.53	1/1189 (0.1%)	0.92	9/1609 (0.6%)
1	K	0.42	1/1228 (0.1%)	0.88	5/1661 (0.3%)
1	L	0.37	0/1118	0.87	3/1511 (0.2%)
All	All	0.42	7/14063 (0.0%)	0.87	51/19004 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	2	LYS	C-N	10.66	1.48	1.33
1	I	25	SER	C-N	-7.73	1.22	1.33
1	K	151	MET	C-N	-6.44	1.24	1.33
1	C	153	ASN	C-N	-5.86	1.25	1.33
1	D	151	MET	C-N	-5.79	1.25	1.33
1	H	153	ASN	C-N	-5.69	1.25	1.33
1	F	151	MET	C-N	-5.67	1.25	1.33

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	82	SER	N-CA-C	8.38	123.73	107.98
1	F	82	SER	N-CA-C	8.35	123.69	107.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	82	SER	N-CA-C	8.14	123.29	107.98
1	G	82	SER	N-CA-C	8.08	123.17	107.98
1	L	82	SER	N-CA-C	7.95	121.98	107.99
1	J	19	GLU	CA-C-N	-7.93	109.93	119.84
1	J	19	GLU	C-N-CA	-7.93	109.93	119.84
1	B	82	SER	N-CA-C	7.91	122.85	107.98
1	D	82	SER	N-CA-C	7.91	122.85	107.98
1	A	82	SER	N-CA-C	7.81	121.73	107.99
1	C	82	SER	N-CA-C	7.76	122.58	107.98
1	E	82	SER	N-CA-C	7.72	121.58	107.99
1	K	82	SER	N-CA-C	7.66	122.39	107.98
1	H	82	SER	N-CA-C	7.64	122.34	107.98
1	F	73	ILE	N-CA-C	6.81	117.92	108.12
1	J	20	PRO	N-CA-C	-6.39	99.30	112.47
1	I	73	ILE	N-CA-C	6.37	117.30	108.12
1	B	73	ILE	N-CA-C	6.37	117.29	108.12
1	G	73	ILE	N-CA-C	6.32	117.22	108.12
1	J	73	ILE	N-CA-C	6.20	117.05	108.12
1	C	73	ILE	N-CA-C	6.17	117.01	108.12
1	F	121	LYS	N-CA-C	-6.02	104.80	111.36
1	K	73	ILE	N-CA-C	6.00	116.77	108.12
1	J	21	HIS	N-CA-C	5.84	118.93	110.28
1	I	117	SER	N-CA-C	5.84	118.11	111.11
1	D	73	ILE	N-CA-C	5.82	116.50	108.12
1	H	73	ILE	N-CA-C	5.82	116.49	108.12
1	E	121	LYS	N-CA-C	-5.76	105.09	111.36
1	B	117	SER	N-CA-C	5.74	118.00	111.11
1	A	73	ILE	N-CA-C	5.72	116.53	108.17
1	E	73	ILE	N-CA-C	5.71	116.35	108.12
1	G	121	LYS	N-CA-C	-5.67	105.19	111.36
1	L	73	ILE	N-CA-C	5.65	116.41	108.17
1	L	117	SER	N-CA-C	5.63	118.14	111.33
1	A	117	SER	N-CA-C	5.59	118.10	111.33
1	C	121	LYS	N-CA-C	-5.50	105.36	111.36
1	A	121	LYS	N-CA-C	-5.45	105.42	111.36
1	K	19	GLU	CA-C-N	-5.36	113.64	120.23
1	K	19	GLU	C-N-CA	-5.36	113.64	120.23
1	J	2	LYS	O-C-N	5.30	129.64	122.59
1	J	121	LYS	N-CA-C	-5.23	105.66	111.36
1	C	83	VAL	N-CA-C	-5.21	104.57	111.09
1	D	121	LYS	N-CA-C	-5.21	105.68	111.36
1	K	121	LYS	N-CA-C	-5.21	105.68	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	117	SER	N-CA-C	5.20	117.63	111.33
1	F	117	SER	N-CA-C	5.19	117.34	111.11
1	C	117	SER	N-CA-C	5.19	117.61	111.33
1	I	121	LYS	N-CA-C	-5.18	105.72	111.36
1	J	83	VAL	N-CA-C	-5.16	104.64	111.09
1	H	74	ASN	CA-C-N	5.13	125.21	119.87
1	H	74	ASN	C-N-CA	5.13	125.21	119.87

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1156	0	1141	43	0
1	B	1089	0	1079	28	0
1	C	1161	0	1149	29	0
1	D	1195	0	1178	38	0
1	E	1130	0	1120	27	0
1	F	1145	0	1134	37	0
1	G	1106	0	1097	34	0
1	H	1161	0	1149	29	0
1	I	1115	0	1103	48	0
1	J	1156	0	1141	22	0
1	K	1195	0	1178	33	0
1	L	1089	0	1079	31	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	15	0	0	0	0
2	D	15	0	0	0	0
2	E	15	0	0	0	0
2	F	20	0	0	0	0
2	G	5	0	0	0	0
2	H	20	0	0	0	0
2	I	5	0	0	0	0
2	J	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	K	15	0	0	0	0
2	L	10	0	0	0	0
3	B	8	0	12	0	0
3	E	8	0	12	0	0
3	H	8	0	12	0	0
3	K	8	0	12	0	0
4	A	195	0	0	8	0
4	B	204	0	0	7	0
4	C	216	0	0	4	0
4	D	203	0	0	5	0
4	E	162	0	0	1	0
4	F	216	0	0	0	0
4	G	156	0	0	3	0
4	H	199	0	0	6	0
4	I	179	0	0	14	0
4	J	152	0	0	3	0
4	K	199	0	0	3	0
4	L	181	0	0	5	0
All	All	16147	0	13596	358	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 13.

All (358) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:125[B]:CYS:SG	1:F:125[B]:CYS:CB	2.30	1.19
1:F:125[C]:CYS:CB	1:F:125[C]:CYS:SG	2.32	1.17
1:I:79:THR:HG21	1:I:100:HIS:HE1	1.10	1.10
1:D:19:GLU:HB3	1:D:20:PRO:HD3	1.35	1.07
1:G:75:PRO:HG2	1:G:79:THR:CG2	1.89	1.02
1:G:75:PRO:CG	1:G:79:THR:HG22	1.89	1.01
1:I:79:THR:HG21	1:I:100:HIS:CE1	1.96	0.99
1:I:75:PRO:HG2	1:I:79:THR:HG22	1.02	0.99
1:B:125[A]:CYS:SG	1:H:123:VAL:HB	2.03	0.99
1:A:125[A]:CYS:SG	1:D:123:VAL:HB	2.04	0.98
1:G:79:THR:HG21	1:G:100:HIS:CE1	1.98	0.97
1:G:76:GLY:O	1:G:79:THR:HG23	1.66	0.95
1:G:75:PRO:HG2	1:G:79:THR:HG22	0.95	0.93
1:F:123:VAL:HB	1:I:125[A]:CYS:SG	2.09	0.92
1:C:125[A]:CYS:SG	1:L:123:VAL:HB	2.09	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:76:GLY:O	1:I:79:THR:HG23	1.69	0.92
1:L:103:ASN:HA	1:L:125[B]:CYS:SG	2.10	0.92
1:K:19:GLU:HB2	1:K:20:PRO:HD3	1.52	0.90
1:F:125[A]:CYS:SG	1:I:123:VAL:HB	2.12	0.90
1:A:75:PRO:HG2	1:A:79:THR:OG1	1.71	0.89
1:G:103:ASN:HA	1:G:125[B]:CYS:SG	2.12	0.89
1:A:103:ASN:HA	1:A:125[B]:CYS:SG	2.12	0.89
1:C:123:VAL:HB	1:L:125[A]:CYS:SG	2.13	0.89
1:I:75:PRO:HG2	1:I:79:THR:CG2	1.98	0.87
1:J:19:GLU:OE2	1:J:19:GLU:HA	1.73	0.87
1:E:123:VAL:HB	1:K:125[A]:CYS:SG	2.14	0.86
1:I:75:PRO:CG	1:I:79:THR:HG22	1.98	0.86
1:G:79:THR:HG21	1:G:100:HIS:HE1	1.36	0.85
1:B:103:ASN:HA	1:B:125[B]:CYS:SG	2.17	0.85
1:G:125[A]:CYS:SG	1:J:123:VAL:HB	2.17	0.83
1:J:21:HIS:HB3	4:J:2152:HOH:O	1.77	0.83
1:K:103:ASN:HA	1:K:125[C]:CYS:SG	2.19	0.82
1:C:103:ASN:HA	1:C:125[C]:CYS:SG	2.20	0.81
1:J:102:SER:O	1:J:125[C]:CYS:SG	2.39	0.80
1:E:102:SER:O	1:E:125[B]:CYS:SG	2.40	0.79
1:E:125[A]:CYS:SG	1:K:123:VAL:HB	2.23	0.79
1:G:123:VAL:HB	1:J:125[A]:CYS:SG	2.22	0.79
1:I:103:ASN:HA	1:I:125[B]:CYS:SG	2.23	0.79
1:L:102:SER:O	1:L:125[B]:CYS:SG	2.41	0.79
1:A:2:LYS:HD3	1:A:143:LEU:HB3	1.65	0.79
1:J:104:VAL:HG13	1:J:125[C]:CYS:SG	2.24	0.76
1:F:102:SER:O	1:F:125[B]:CYS:SG	2.43	0.76
1:J:19:GLU:O	1:J:21:HIS:N	2.18	0.76
1:A:123:VAL:HB	1:D:125[A]:CYS:SG	2.24	0.76
1:D:102:SER:O	1:D:125[B]:CYS:SG	2.44	0.76
1:B:102:SER:O	1:B:125[B]:CYS:SG	2.44	0.75
1:D:19:GLU:HB3	1:D:20:PRO:CD	2.16	0.75
1:A:123:VAL:HG12	1:A:125[C]:CYS:SG	2.27	0.75
1:H:102:SER:O	1:H:125[A]:CYS:SG	2.44	0.74
1:K:102:SER:O	1:K:125[C]:CYS:SG	2.45	0.74
1:L:100:HIS:HD2	1:L:125[C]:CYS:SG	2.10	0.74
1:F:103:ASN:HA	1:F:125[B]:CYS:SG	2.28	0.74
1:A:100:HIS:HD2	1:A:125[C]:CYS:SG	2.11	0.73
1:D:103:ASN:HA	1:D:125[B]:CYS:SG	2.28	0.73
1:A:40:GLN:HG3	4:A:2050:HOH:O	1.89	0.73
1:J:100:HIS:HB2	1:J:125[C]:CYS:SG	2.29	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:145:LYS:HE3	4:I:2171:HOH:O	1.89	0.73
1:A:18:ARG:H	1:A:18:ARG:HD3	1.52	0.72
1:I:102:SER:O	1:I:125[B]:CYS:SG	2.46	0.72
1:G:102:SER:O	1:G:125[B]:CYS:SG	2.48	0.72
1:I:100:HIS:HD2	1:I:125[C]:CYS:SG	2.12	0.72
1:A:18:ARG:H	1:A:18:ARG:CD	2.03	0.72
1:I:145:LYS:HG3	4:I:2178:HOH:O	1.89	0.71
1:B:100:HIS:HD2	1:B:125[C]:CYS:SG	2.13	0.71
1:L:144:GLN:C	1:L:146:ILE:H	1.97	0.71
1:B:123:VAL:HG12	1:B:125[C]:CYS:SG	2.31	0.70
1:F:123:VAL:HG12	1:F:125[C]:CYS:SG	2.32	0.70
1:J:123:VAL:HG12	1:J:125[B]:CYS:SG	2.31	0.70
1:C:102:SER:O	1:C:125[C]:CYS:SG	2.50	0.69
1:D:36:GLN:O	1:D:40:GLN:HG3	1.93	0.68
1:F:59:ASN:O	1:F:63:GLN:HG3	1.94	0.68
1:G:3:LYS:HD2	1:G:47:ASP:OD2	1.94	0.68
1:A:18:ARG:HD3	1:A:18:ARG:N	2.08	0.68
1:J:59:ASN:O	1:J:63:GLN:HG3	1.92	0.68
1:F:100:HIS:HD2	1:F:125[C]:CYS:SG	2.16	0.67
1:A:100:HIS:ND1	4:A:2136:HOH:O	2.26	0.67
1:H:123:VAL:HG12	1:H:125[B]:CYS:SG	2.35	0.67
1:I:123:VAL:HG12	1:I:125[C]:CYS:SG	2.35	0.67
1:A:102:SER:O	1:A:125[B]:CYS:SG	2.52	0.66
1:F:2:LYS:HG3	1:F:143:LEU:HD13	1.76	0.66
1:F:2:LYS:HE3	1:F:147:GLN:HG2	1.79	0.65
1:L:100:HIS:CD2	1:L:125[C]:CYS:SG	2.90	0.65
1:A:21:HIS:HE1	1:A:30:ASP:OD2	1.80	0.65
1:D:59:ASN:O	1:D:63:GLN:HG3	1.97	0.64
1:G:75:PRO:HG3	1:G:116:LEU:HD12	1.79	0.64
1:J:1:MET:O	1:J:2:LYS:O	2.16	0.64
1:C:35:LEU:HD11	1:C:132:TYR:HB3	1.79	0.64
1:B:2:LYS:HB3	1:B:44:TYR:CD1	2.33	0.64
1:K:19:GLU:CB	1:K:20:PRO:HD3	2.26	0.63
1:A:17:LYS:HE2	4:A:2021:HOH:O	2.00	0.62
1:F:123:VAL:HB	1:I:125[A]:CYS:HG	1.64	0.62
1:I:140:ILE:O	1:I:144:GLN:HG3	2.00	0.62
1:C:146:ILE:O	1:C:150:GLU:HG3	2.00	0.61
1:C:104:VAL:N	1:C:125[C]:CYS:SG	2.70	0.61
1:H:2:LYS:HE2	4:H:2077:HOH:O	2.00	0.61
1:I:146:ILE:HD11	4:I:2079:HOH:O	2.00	0.61
1:K:19:GLU:HB2	1:K:20:PRO:CD	2.28	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:34:HIS:HD2	4:L:2026:HOH:O	1.83	0.61
1:L:144:GLN:C	1:L:146:ILE:N	2.55	0.61
1:K:149:GLY:C	1:K:151:MET:H	2.08	0.60
1:F:123:VAL:CB	1:I:125[A]:CYS:SG	2.88	0.60
1:L:121:LYS:HE2	1:L:142:GLU:OE2	2.01	0.60
1:I:75:PRO:HG3	1:I:116:LEU:HD12	1.81	0.60
1:K:104:VAL:N	1:K:125[C]:CYS:SG	2.72	0.60
1:A:17:LYS:HE3	4:B:2140:HOH:O	2.01	0.59
1:E:103:ASN:HA	1:E:125[B]:CYS:SG	2.42	0.59
1:A:3:LYS:HE2	1:A:45:GLU:OE2	2.01	0.59
1:E:123:VAL:HG12	1:E:125[C]:CYS:SG	2.43	0.59
1:A:143:LEU:O	1:A:146:ILE:HG12	2.02	0.59
1:A:100:HIS:CD2	1:A:125[C]:CYS:SG	2.96	0.59
1:K:103:ASN:C	1:K:103:ASN:HD22	2.10	0.59
1:G:100:HIS:HD2	1:G:125[C]:CYS:SG	2.25	0.58
1:C:40:GLN:NE2	4:C:2071:HOH:O	2.35	0.58
1:D:15:LEU:HD11	1:D:101:LEU:HD11	1.84	0.58
1:B:140:ILE:O	1:B:144:GLN:HG3	2.03	0.58
1:D:104:VAL:N	1:D:125[B]:CYS:SG	2.72	0.58
1:B:125[A]:CYS:SG	1:H:123:VAL:CB	2.87	0.58
1:B:104:VAL:N	1:B:125[B]:CYS:SG	2.74	0.58
1:I:37:GLN:HG2	4:I:2055:HOH:O	2.03	0.58
1:I:107:ARG:C	4:I:2126:HOH:O	2.48	0.57
1:C:148:LEU:HG	1:C:152:MET:HE2	1.85	0.57
1:D:123:VAL:HG12	1:D:125[C]:CYS:SG	2.45	0.56
1:E:75:PRO:HG2	1:E:79:THR:HB	1.87	0.56
1:E:53:GLY:HA3	1:F:55:GLU:HB2	1.87	0.56
1:I:25:SER:O	1:I:26:GLN:O	2.24	0.56
1:I:104:VAL:O	4:I:2126:HOH:O	2.18	0.56
1:C:3:LYS:HD2	1:C:47:ASP:OD2	2.06	0.56
1:J:19:GLU:OE2	1:J:19:GLU:CA	2.51	0.56
1:I:25:SER:N	4:I:2031:HOH:O	2.38	0.56
1:G:100:HIS:HB2	1:G:125[C]:CYS:SG	2.45	0.55
1:B:1:MET:HG2	4:B:2002:HOH:O	2.06	0.55
1:G:100:HIS:CD2	1:G:125[C]:CYS:SG	2.99	0.55
1:L:36:GLN:O	1:L:40:GLN:HG3	2.07	0.55
1:L:140:ILE:O	1:L:144:GLN:HG3	2.06	0.55
1:F:2:LYS:HG2	1:F:143:LEU:HB3	1.88	0.55
1:D:75:PRO:HG2	1:D:79:THR:HB	1.87	0.55
1:F:125[A]:CYS:SG	1:I:123:VAL:CB	2.92	0.55
1:I:112:HIS:CD2	4:I:2126:HOH:O	2.60	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:VAL:HB	1:H:125[C]:CYS:SG	2.47	0.55
1:H:100:HIS:HB2	1:H:125[A]:CYS:SG	2.46	0.54
1:B:52:ASN:HD22	1:C:59:ASN:HD21	1.55	0.54
1:F:75:PRO:HG2	1:F:116:LEU:HD12	1.90	0.54
1:K:151:MET:HE3	1:K:151:MET:HA	1.89	0.54
1:G:104:VAL:N	1:G:125[B]:CYS:SG	2.78	0.54
1:B:40:GLN:HG3	4:B:2083:HOH:O	2.08	0.54
1:I:18:ARG:O	1:I:19:GLU:C	2.50	0.54
1:H:104:VAL:HG12	4:H:2140:HOH:O	2.08	0.54
1:B:53:GLY:HA3	1:C:55:GLU:HB2	1.89	0.53
1:C:75:PRO:HG2	1:C:79:THR:HB	1.89	0.53
1:H:36:GLN:O	1:H:40:GLN:HG3	2.07	0.53
1:F:104:VAL:N	1:F:125[B]:CYS:SG	2.72	0.53
1:D:45:GLU:HB3	4:D:2083:HOH:O	2.08	0.53
1:A:18:ARG:CD	1:A:18:ARG:N	2.69	0.53
1:I:79:THR:CG2	1:I:100:HIS:HE1	2.01	0.53
1:L:124:ILE:O	1:L:125[C]:CYS:SG	2.67	0.53
1:G:36:GLN:O	1:G:40:GLN:HG3	2.09	0.53
1:H:148:LEU:HG	1:H:152:MET:HE2	1.91	0.53
1:E:140:ILE:O	1:E:144:GLN:HG3	2.09	0.53
1:K:149:GLY:C	1:K:151:MET:N	2.66	0.53
1:B:28:LEU:HD23	4:B:2008:HOH:O	2.08	0.52
1:C:55:GLU:CD	4:C:2105:HOH:O	2.51	0.52
1:E:104:VAL:HG13	1:E:125[B]:CYS:SG	2.49	0.52
1:E:104:VAL:N	1:E:125[B]:CYS:SG	2.74	0.52
1:E:140:ILE:HG22	1:E:144:GLN:HE21	1.75	0.52
1:G:33:GLN:O	1:G:37:GLN:HG3	2.10	0.52
1:H:2:LYS:HE2	4:H:2178:HOH:O	2.10	0.52
1:A:2:LYS:CD	1:A:143:LEU:HB3	2.35	0.52
1:F:1:MET:HG3	1:F:2:LYS:H	1.75	0.52
1:I:2:LYS:HE3	4:I:2175:HOH:O	2.10	0.52
1:I:104:VAL:N	1:I:125[B]:CYS:SG	2.73	0.52
1:F:2:LYS:HE3	1:F:147:GLN:CG	2.40	0.51
1:H:53:GLY:HA3	1:I:55:GLU:HB2	1.93	0.51
1:B:100:HIS:CD2	1:B:125[C]:CYS:SG	3.01	0.51
1:D:19:GLU:O	1:D:21:HIS:N	2.41	0.51
1:F:143:LEU:O	1:F:147:GLN:HG3	2.11	0.51
1:I:100:HIS:CD2	1:I:125[C]:CYS:SG	2.99	0.51
1:K:148:LEU:C	1:K:148:LEU:HD23	2.35	0.51
1:I:25:SER:O	1:I:26:GLN:C	2.50	0.51
1:B:125[A]:CYS:HG	1:H:123:VAL:HB	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:116:LEU:HG	4:G:2094:HOH:O	2.10	0.51
1:A:2:LYS:HE3	1:A:143:LEU:O	2.11	0.51
1:A:125[A]:CYS:SG	1:D:123:VAL:CB	2.89	0.51
1:E:1:MET:HE3	1:E:45:GLU:HG3	1.92	0.51
1:I:141:SER:O	1:I:145:LYS:HG2	2.11	0.51
1:I:26:GLN:HG2	4:I:2034:HOH:O	2.11	0.51
1:K:75:PRO:HG2	1:K:79:THR:HB	1.92	0.51
1:J:123:VAL:CG1	1:J:125[B]:CYS:SG	2.99	0.50
1:A:75:PRO:HG2	1:A:79:THR:HG1	1.76	0.50
1:B:130:LYS:HG2	4:B:2189:HOH:O	2.11	0.50
1:C:125[A]:CYS:SG	1:L:123:VAL:CB	2.93	0.50
1:F:3:LYS:HD2	1:F:47:ASP:OD2	2.12	0.50
1:I:40:GLN:HG3	4:I:2073:HOH:O	2.11	0.50
1:B:3:LYS:HD3	4:B:2125:HOH:O	2.10	0.50
1:H:100:HIS:HD2	1:H:125[B]:CYS:SG	2.35	0.50
1:E:2:LYS:HD3	1:E:143:LEU:HB3	1.94	0.50
1:D:65:PHE:CE2	1:F:18:ARG:HG3	2.47	0.50
1:I:130:LYS:HE3	1:I:134:TYR:OH	2.11	0.50
1:K:19:GLU:HG2	4:L:2105:HOH:O	2.11	0.50
1:A:123:VAL:CG1	1:A:125[C]:CYS:SG	2.99	0.49
1:I:123:VAL:CG1	1:I:125[C]:CYS:SG	3.00	0.49
1:J:55:GLU:HB2	1:L:53:GLY:HA3	1.93	0.49
1:E:3:LYS:HD3	1:E:47:ASP:OD2	2.12	0.49
1:K:53:GLY:HA3	1:L:55:GLU:HB2	1.93	0.49
1:A:2:LYS:HZ2	1:A:146:ILE:HD11	1.78	0.49
1:D:3:LYS:HD2	1:D:47:ASP:OD2	2.11	0.49
1:K:145:LYS:HE2	4:K:2186:HOH:O	2.11	0.49
1:H:97:ILE:HD11	1:H:138:PHE:CD2	2.48	0.49
1:J:28:LEU:HD21	1:J:50:GLN:HG3	1.94	0.49
1:J:53:GLY:HA3	1:K:55:GLU:HB2	1.94	0.49
1:A:55:GLU:HB2	1:C:53:GLY:HA3	1.94	0.49
1:E:33:GLN:O	1:E:37:GLN:HG3	2.13	0.49
1:J:75:PRO:HG2	1:J:79:THR:HB	1.95	0.49
1:H:123:VAL:CG1	1:H:125[B]:CYS:SG	3.01	0.49
1:A:123:VAL:CB	1:D:125[A]:CYS:SG	2.98	0.48
1:I:33:GLN:NE2	4:I:2047:HOH:O	2.44	0.48
1:E:123:VAL:CB	1:K:125[A]:CYS:SG	2.97	0.48
1:E:59:ASN:O	1:E:63:GLN:HG3	2.13	0.48
1:I:3:LYS:HG3	1:I:45:GLU:HG3	1.94	0.48
1:L:26:GLN:N	4:L:2029:HOH:O	2.45	0.48
1:G:130:LYS:HG3	4:G:2143:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:LYS:HD2	1:B:67:ASN:O	2.14	0.48
1:B:1:MET:O	1:B:2:LYS:C	2.56	0.48
1:G:1:MET:SD	1:G:1:MET:C	2.97	0.48
1:C:97:ILE:HD11	1:C:138:PHE:CD2	2.48	0.48
1:D:19:GLU:HA	4:D:2023:HOH:O	2.13	0.48
1:F:123:VAL:CG1	1:F:125[C]:CYS:SG	3.00	0.47
1:H:147:GLN:HG3	4:H:2178:HOH:O	2.14	0.47
1:K:46:LEU:HD23	1:K:46:LEU:C	2.38	0.47
1:B:52:ASN:HD22	1:C:59:ASN:ND2	2.11	0.47
1:D:55:GLU:HB2	1:F:53:GLY:HA3	1.95	0.47
1:I:97:ILE:HD11	1:I:138:PHE:CD2	2.49	0.47
1:C:130:LYS:HG3	4:C:2185:HOH:O	2.14	0.47
1:A:37:GLN:HG2	4:A:2048:HOH:O	2.15	0.47
1:E:46:LEU:C	1:E:46:LEU:HD23	2.40	0.47
1:K:15:LEU:HD21	1:K:101:LEU:CD1	2.44	0.47
1:C:33:GLN:O	1:C:37:GLN:HG3	2.15	0.47
1:L:1:MET:CB	4:L:2067:HOH:O	2.62	0.47
1:G:104:VAL:HG12	4:G:2115:HOH:O	2.15	0.47
1:H:146:ILE:O	1:H:150:GLU:HG3	2.15	0.47
1:D:53:GLY:HA3	1:E:55:GLU:HB2	1.96	0.46
1:I:2:LYS:O	1:I:45:GLU:HG2	2.15	0.46
1:L:103:ASN:CA	1:L:125[B]:CYS:SG	2.94	0.46
1:C:123:VAL:HG12	1:C:125[B]:CYS:SG	2.55	0.46
1:F:147:GLN:O	1:F:150:GLU:HG2	2.15	0.46
1:H:104:VAL:HG13	1:H:125[A]:CYS:SG	2.55	0.46
1:A:53:GLY:HA3	1:B:55:GLU:HB2	1.97	0.46
1:D:46:LEU:HD23	1:D:46:LEU:C	2.41	0.46
1:B:103:ASN:CA	1:B:125[B]:CYS:SG	2.97	0.46
1:G:55:GLU:HB2	1:I:53:GLY:HA3	1.98	0.46
1:G:97:ILE:HD11	1:G:138:PHE:CD2	2.51	0.46
1:L:75:PRO:HG2	1:L:116:LEU:HD12	1.99	0.45
1:F:1:MET:CG	1:F:2:LYS:N	2.77	0.45
1:J:100:HIS:CB	1:J:125[C]:CYS:SG	3.03	0.45
1:L:97:ILE:HD13	1:L:135:ALA:HA	1.98	0.45
1:G:130:LYS:HE3	1:G:134:TYR:OH	2.16	0.45
1:K:28:LEU:HD21	1:K:50:GLN:HG3	1.98	0.45
1:G:103:ASN:CA	1:G:125[B]:CYS:SG	2.95	0.45
1:L:144:GLN:O	1:L:146:ILE:N	2.49	0.45
1:D:130:LYS:HG2	4:D:2182:HOH:O	2.17	0.45
1:G:46:LEU:C	1:G:46:LEU:HD23	2.41	0.45
1:B:1:MET:O	1:B:3:LYS:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:1:MET:HG3	1:F:2:LYS:N	2.31	0.44
1:L:79:THR:OG1	1:L:100:HIS:HE1	2.00	0.44
1:L:104:VAL:HG12	4:L:2141:HOH:O	2.17	0.44
1:A:21:HIS:CE1	1:A:30:ASP:OD2	2.66	0.44
1:D:97:ILE:HD11	1:D:138:PHE:CD2	2.52	0.44
1:H:130:LYS:HE3	1:H:133:ASP:OD1	2.18	0.44
1:A:36:GLN:NE2	4:A:2048:HOH:O	2.44	0.44
1:A:47:ASP:HB3	4:A:2106:HOH:O	2.17	0.44
1:A:141:SER:O	1:A:145:LYS:HG3	2.18	0.44
1:D:18:ARG:NH2	4:D:2020:HOH:O	2.50	0.44
1:C:100:HIS:HD2	1:C:125[B]:CYS:SG	2.41	0.44
1:J:28:LEU:CD2	1:J:50:GLN:HG3	2.48	0.44
1:D:28:LEU:O	1:D:32:GLU:HG3	2.18	0.43
1:G:53:GLY:HA3	1:H:55:GLU:HB2	1.99	0.43
1:H:79:THR:HG23	1:H:80:HIS:CE1	2.53	0.43
1:L:104:VAL:N	1:L:125[B]:CYS:SG	2.86	0.43
1:G:79:THR:CG2	1:G:100:HIS:HE1	2.18	0.43
1:C:10:PRO:HA	1:C:52:ASN:OD1	2.19	0.43
1:C:103:ASN:CA	1:C:125[C]:CYS:SG	2.99	0.43
4:C:2067:HOH:O	1:L:112:HIS:HD2	2.00	0.43
1:D:91:ALA:O	1:F:18:ARG:HD3	2.18	0.43
1:E:125[A]:CYS:SG	1:K:123:VAL:CB	3.02	0.43
1:H:121:LYS:NZ	4:H:2169:HOH:O	2.51	0.43
1:K:103:ASN:CA	1:K:125[C]:CYS:SG	2.99	0.43
1:E:130:LYS:HE3	1:E:134:TYR:OH	2.17	0.43
1:H:140:ILE:O	1:H:144:GLN:HG3	2.18	0.43
1:H:143:LEU:O	1:H:147:GLN:HG3	2.18	0.43
1:E:123:VAL:CG1	1:E:125[C]:CYS:SG	3.07	0.43
1:K:10:PRO:HA	1:K:52:ASN:OD1	2.18	0.43
1:L:124:ILE:C	1:L:125[C]:CYS:SG	3.02	0.43
1:A:18:ARG:O	1:A:19:GLU:C	2.61	0.43
1:G:3:LYS:HE2	1:G:67:ASN:O	2.19	0.43
1:G:139:ALA:O	1:G:143:LEU:HG	2.19	0.43
1:H:46:LEU:C	1:H:46:LEU:HD23	2.45	0.42
1:K:151:MET:HA	1:K:151:MET:CE	2.46	0.42
1:E:10:PRO:HA	1:E:52:ASN:OD1	2.20	0.42
1:L:123:VAL:HG12	1:L:125[C]:CYS:SG	2.58	0.42
1:A:103:ASN:CA	1:A:125[B]:CYS:SG	2.96	0.42
1:E:23:TYR:CE2	1:E:101:LEU:HB3	2.53	0.42
1:K:130:LYS:HG3	4:K:2171:HOH:O	2.19	0.42
1:C:3:LYS:HE2	1:C:67:ASN:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:75:PRO:HG2	1:H:116:LEU:HD12	2.01	0.42
1:K:33:GLN:O	1:K:37:GLN:HG3	2.19	0.42
1:A:46:LEU:C	1:A:46:LEU:HD23	2.45	0.42
1:A:35:LEU:HD11	1:A:132:TYR:HB3	2.01	0.42
1:C:75:PRO:HG3	1:C:116:LEU:HD12	2.02	0.42
1:D:35:LEU:HG	1:D:136:LEU:HD22	2.02	0.42
1:D:100:HIS:HD2	1:D:125[C]:CYS:SG	2.43	0.42
1:L:72:ILE:HD11	1:L:136:LEU:HA	2.02	0.42
1:D:4:ILE:O	1:D:46:LEU:HA	2.20	0.42
1:A:104:VAL:HG12	4:A:2146:HOH:O	2.19	0.41
1:B:26:GLN:N	4:B:2039:HOH:O	2.52	0.41
1:D:28:LEU:CD2	1:D:50:GLN:HG3	2.50	0.41
1:F:130:LYS:HE3	1:I:141:SER:CB	2.50	0.41
1:L:46:LEU:C	1:L:46:LEU:HD23	2.45	0.41
1:G:28:LEU:HD21	1:G:50:GLN:HG3	2.03	0.41
1:E:2:LYS:CD	1:E:143:LEU:HB3	2.49	0.41
1:J:28:LEU:HD23	4:J:2002:HOH:O	2.21	0.41
1:A:104:VAL:N	1:A:125[B]:CYS:SG	2.83	0.41
1:F:1:MET:CG	1:F:2:LYS:H	2.33	0.41
1:C:127:LEU:HD22	1:L:138:PHE:CD2	2.56	0.41
1:D:3:LYS:HE2	1:D:67:ASN:O	2.21	0.41
1:K:4:ILE:O	1:K:46:LEU:HA	2.20	0.41
1:K:103:ASN:C	1:K:103:ASN:ND2	2.79	0.41
1:A:79:THR:OG1	1:A:100:HIS:HE1	2.03	0.41
1:D:65:PHE:CZ	1:F:18:ARG:HG3	2.56	0.41
1:A:17:LYS:HB3	1:A:18:ARG:H	1.68	0.41
1:B:123:VAL:CG1	1:B:125[C]:CYS:SG	3.03	0.41
1:C:153:ASN:HD22	1:C:153:ASN:HA	1.64	0.41
1:D:103:ASN:CA	1:D:125[B]:CYS:SG	3.06	0.41
1:D:121:LYS:NZ	4:D:2174:HOH:O	2.54	0.41
1:H:2:LYS:HG2	4:H:2186:HOH:O	2.20	0.41
1:J:18:ARG:NH1	4:J:2024:HOH:O	2.46	0.41
1:K:104:VAL:HG12	4:K:2131:HOH:O	2.21	0.41
1:F:2:LYS:HD3	1:F:69:ASP:CG	2.47	0.40
1:G:15:LEU:HD11	1:G:101:LEU:HD11	2.03	0.40
1:I:103:ASN:CA	1:I:125[B]:CYS:SG	3.02	0.40
1:I:26:GLN:N	4:I:2037:HOH:O	2.39	0.40
1:A:121:LYS:NZ	4:A:2178:HOH:O	2.51	0.40
1:E:146:ILE:HG13	4:E:2158:HOH:O	2.20	0.40
1:K:103:ASN:ND2	1:K:105:HIS:H	2.20	0.40
1:D:104:VAL:HG13	1:D:125[B]:CYS:SG	2.61	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:123:VAL:CG1	1:D:125[C]:CYS:SG	3.09	0.40
1:F:75:PRO:CG	1:F:116:LEU:HD12	2.52	0.40
1:I:104:VAL:HG12	4:I:2124:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	146/154 (95%)	139 (95%)	7 (5%)	0	100	100
1	B	136/154 (88%)	133 (98%)	2 (2%)	1 (1%)	18	7
1	C	146/154 (95%)	144 (99%)	2 (1%)	0	100	100
1	D	152/154 (99%)	146 (96%)	5 (3%)	1 (1%)	18	7
1	E	141/154 (92%)	139 (99%)	2 (1%)	0	100	100
1	F	144/154 (94%)	142 (99%)	2 (1%)	0	100	100
1	G	138/154 (90%)	136 (99%)	2 (1%)	0	100	100
1	H	146/154 (95%)	144 (99%)	2 (1%)	0	100	100
1	I	139/154 (90%)	135 (97%)	3 (2%)	1 (1%)	18	7
1	J	146/154 (95%)	142 (97%)	3 (2%)	1 (1%)	18	7
1	K	152/154 (99%)	147 (97%)	3 (2%)	2 (1%)	9	2
1	L	136/154 (88%)	133 (98%)	3 (2%)	0	100	100
All	All	1722/1848 (93%)	1680 (98%)	36 (2%)	6 (0%)	30	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	26	GLN

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Mol	Chain	Res	Type
1	D	21	HIS
1	J	2	LYS
1	K	151	MET
1	K	19	GLU
1	B	2	LYS

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	125/129 (97%)	123 (98%)	2 (2%)	55	41
1	B	118/129 (92%)	117 (99%)	1 (1%)	73	65
1	C	126/129 (98%)	126 (100%)	0	100	100
1	D	129/129 (100%)	126 (98%)	3 (2%)	44	27
1	E	122/129 (95%)	122 (100%)	0	100	100
1	F	124/129 (96%)	124 (100%)	0	100	100
1	G	120/129 (93%)	118 (98%)	2 (2%)	53	38
1	H	126/129 (98%)	126 (100%)	0	100	100
1	I	121/129 (94%)	119 (98%)	2 (2%)	53	38
1	J	125/129 (97%)	122 (98%)	3 (2%)	43	26
1	K	129/129 (100%)	128 (99%)	1 (1%)	73	65
1	L	118/129 (92%)	117 (99%)	1 (1%)	73	65
All	All	1483/1548 (96%)	1468 (99%)	15 (1%)	68	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	22	ILE
1	B	1	MET
1	D	19	GLU
1	D	21	HIS

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Mol	Chain	Res	Type
1	D	35	LEU
1	G	1	MET
1	G	35	LEU
1	I	26	GLN
1	I	79	THR
1	J	19	GLU
1	J	21	HIS
1	J	35	LEU
1	K	151	MET
1	L	1	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (74) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	26	GLN
1	A	36	GLN
1	A	37	GLN
1	A	40	GLN
1	A	42	GLN
1	A	100	HIS
1	A	105	HIS
1	A	144	GLN
1	B	26	GLN
1	B	40	GLN
1	B	100	HIS
1	B	144	GLN
1	C	33	GLN
1	C	37	GLN
1	C	40	GLN
1	C	42	GLN
1	C	59	ASN
1	C	63	GLN
1	C	112	HIS
1	C	144	GLN
1	C	153	ASN
1	D	21	HIS
1	D	26	GLN
1	D	33	GLN
1	D	34	HIS
1	D	66	GLN
1	D	67	ASN

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Mol	Chain	Res	Type
1	D	105	HIS
1	D	144	GLN
1	E	26	GLN
1	E	40	GLN
1	E	63	GLN
1	E	144	GLN
1	F	26	GLN
1	F	42	GLN
1	F	67	ASN
1	F	100	HIS
1	F	105	HIS
1	F	144	GLN
1	G	33	GLN
1	G	36	GLN
1	G	100	HIS
1	G	105	HIS
1	G	144	GLN
1	H	13	ASN
1	H	33	GLN
1	H	34	HIS
1	H	37	GLN
1	H	66	GLN
1	H	67	ASN
1	H	100	HIS
1	H	144	GLN
1	I	26	GLN
1	I	40	GLN
1	I	66	GLN
1	I	100	HIS
1	I	105	HIS
1	J	36	GLN
1	J	37	GLN
1	J	40	GLN
1	J	42	GLN
1	J	105	HIS
1	J	112	HIS
1	J	144	GLN
1	K	34	HIS
1	K	37	GLN
1	K	40	GLN
1	K	66	GLN
1	K	103	ASN

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Mol	Chain	Res	Type
1	K	144	GLN
1	L	26	GLN
1	L	34	HIS
1	L	100	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

35 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	D	1153	-	4,4,4	0.38	0	6,6,6	0.06	0
2	SO4	F	1155	-	4,4,4	0.37	0	6,6,6	0.14	0
2	SO4	C	1156	-	4,4,4	0.36	0	6,6,6	0.13	0
3	TRS	H	1158	-	7,7,7	1.36	0	9,9,9	1.96	4 (44%)
3	TRS	K	1155	-	7,7,7	1.42	0	9,9,9	1.94	4 (44%)
2	SO4	H	1157	-	4,4,4	0.37	0	6,6,6	0.12	0
2	SO4	J	1147	-	4,4,4	0.35	0	6,6,6	0.08	0
2	SO4	G	1147	-	4,4,4	0.34	0	6,6,6	0.10	0
2	SO4	A	1148	-	4,4,4	0.32	0	6,6,6	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	F	1154	-	4,4,4	0.39	0	6,6,6	0.16	0
2	SO4	B	1148	-	4,4,4	0.39	0	6,6,6	0.06	0
2	SO4	J	1148	-	4,4,4	0.36	0	6,6,6	0.06	0
2	SO4	C	1155	-	4,4,4	0.38	0	6,6,6	0.07	0
2	SO4	A	1147	-	4,4,4	0.36	0	6,6,6	0.14	0
2	SO4	D	1152	-	4,4,4	0.31	0	6,6,6	0.11	0
2	SO4	H	1155	-	4,4,4	0.35	0	6,6,6	0.11	0
2	SO4	E	1147	-	4,4,4	0.34	0	6,6,6	0.10	0
2	SO4	F	1152	-	4,4,4	0.38	0	6,6,6	0.18	0
2	SO4	K	1154	-	4,4,4	0.37	0	6,6,6	0.08	0
2	SO4	E	1149	-	4,4,4	0.37	0	6,6,6	0.11	0
2	SO4	H	1156	-	4,4,4	0.38	0	6,6,6	0.08	0
3	TRS	B	1149	-	7,7,7	1.52	2 (28%)	9,9,9	2.00	4 (44%)
2	SO4	C	1154	-	4,4,4	0.34	0	6,6,6	0.12	0
2	SO4	E	1148	-	4,4,4	0.38	0	6,6,6	0.08	0
3	TRS	E	1150	-	7,7,7	1.49	0	9,9,9	2.07	4 (44%)
2	SO4	K	1153	-	4,4,4	0.38	0	6,6,6	0.09	0
2	SO4	L	1147	-	4,4,4	0.33	0	6,6,6	0.11	0
2	SO4	B	1147	-	4,4,4	0.32	0	6,6,6	0.13	0
2	SO4	I	1147	-	4,4,4	0.37	0	6,6,6	0.11	0
2	SO4	D	1154	-	4,4,4	0.36	0	6,6,6	0.11	0
2	SO4	F	1153	-	4,4,4	0.37	0	6,6,6	0.10	0
2	SO4	J	1149	-	4,4,4	0.38	0	6,6,6	0.13	0
2	SO4	K	1152	-	4,4,4	0.34	0	6,6,6	0.08	0
2	SO4	L	1148	-	4,4,4	0.36	0	6,6,6	0.13	0
2	SO4	H	1154	-	4,4,4	0.43	0	6,6,6	0.18	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	TRS	E	1150	-	-	0/9/9/9	-
3	TRS	K	1155	-	-	0/9/9/9	-
3	TRS	B	1149	-	-	0/9/9/9	-
3	TRS	H	1158	-	-	0/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1149	TRS	C-N	-2.16	1.42	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1149	TRS	C1-C	2.05	1.58	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1149	TRS	C2-C-N	3.10	116.07	108.17
3	K	1155	TRS	C2-C-N	3.04	115.91	108.17
3	E	1150	TRS	C2-C-N	3.00	115.82	108.17
3	H	1158	TRS	C2-C-N	2.99	115.81	108.17
3	B	1149	TRS	O1-C1-C	-2.76	103.20	110.88
3	H	1158	TRS	C2-C-C1	-2.67	103.54	110.66
3	E	1150	TRS	C2-C-C1	-2.67	103.55	110.66
3	E	1150	TRS	O1-C1-C	-2.66	103.46	110.88
3	K	1155	TRS	C2-C-C1	-2.66	103.57	110.66
3	B	1149	TRS	C2-C-C1	-2.61	103.71	110.66
3	E	1150	TRS	C3-C-N	2.56	114.70	108.17
3	B	1149	TRS	C3-C-N	2.39	114.25	108.17
3	K	1155	TRS	O1-C1-C	-2.38	104.24	110.88
3	K	1155	TRS	C3-C-N	2.38	114.24	108.17
3	H	1158	TRS	O1-C1-C	-2.38	104.25	110.88
3	H	1158	TRS	C3-C-N	2.34	114.14	108.17

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(Å²)	Q<0.9
1	A	146/154 (94%)	-0.09	10 (6%)	23	25	11, 18, 49, 66	1 (0%)
1	B	138/154 (89%)	-0.37	3 (2%)	62	66	10, 16, 33, 62	1 (0%)
1	C	148/154 (96%)	-0.21	5 (3%)	48	52	11, 20, 38, 74	1 (0%)
1	D	152/154 (98%)	0.04	11 (7%)	21	23	13, 20, 56, 73	1 (0%)
1	E	143/154 (92%)	0.11	9 (6%)	26	28	14, 23, 48, 73	1 (0%)
1	F	146/154 (94%)	-0.23	7 (4%)	35	39	12, 18, 40, 70	1 (0%)
1	G	140/154 (90%)	0.03	5 (3%)	46	50	13, 22, 45, 69	1 (0%)
1	H	148/154 (96%)	-0.05	5 (3%)	48	52	13, 21, 37, 72	1 (0%)
1	I	141/154 (91%)	-0.20	7 (4%)	34	38	12, 18, 39, 67	1 (0%)
1	J	146/154 (94%)	0.27	11 (7%)	20	22	14, 24, 52, 71	1 (0%)
1	K	152/154 (98%)	0.03	9 (5%)	28	31	15, 21, 50, 74	1 (0%)
1	L	138/154 (89%)	-0.15	4 (2%)	53	58	14, 18, 36, 70	1 (0%)
All	All	1738/1848 (94%)	-0.07	86 (4%)	35	38	10, 20, 45, 74	12 (0%)

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	154	GLY	10.7
1	D	20	PRO	6.8
1	I	25	SER	5.8
1	E	146	ILE	5.6
1	L	146	ILE	5.6
1	A	23	TYR	5.5
1	J	19	GLU	5.4
1	D	21	HIS	5.2
1	J	146	ILE	5.2
1	K	19	GLU	5.1
1	E	22	ILE	5.1

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Mol	Chain	Res	Type	RSRZ
1	C	154	GLY	5.0
1	J	21	HIS	5.0
1	A	146	ILE	5.0
1	J	20	PRO	4.9
1	D	22	ILE	4.9
1	E	1	MET	4.6
1	K	152	MET	4.6
1	F	1	MET	4.5
1	G	1	MET	4.3
1	B	1	MET	4.3
1	K	23	TYR	4.2
1	G	146	ILE	4.2
1	K	151	MET	4.2
1	D	23	TYR	4.2
1	A	24	GLY	4.2
1	J	22	ILE	4.1
1	A	22	ILE	3.9
1	E	23	TYR	3.9
1	E	24	GLY	3.7
1	K	20	PRO	3.7
1	J	1	MET	3.6
1	D	152	MET	3.5
1	B	146	ILE	3.5
1	A	18	ARG	3.5
1	A	20	PRO	3.5
1	E	144	GLN	3.4
1	L	1	MET	3.4
1	F	150	GLU	3.2
1	I	146	ILE	3.2
1	D	24	GLY	3.1
1	I	1	MET	3.1
1	K	21	HIS	3.0
1	J	24	GLY	3.0
1	A	1	MET	3.0
1	F	152	MET	2.9
1	G	18	ARG	2.9
1	F	2	LYS	2.9
1	E	18	ARG	2.8
1	D	18	ARG	2.7
1	A	19	GLU	2.6
1	D	19	GLU	2.6
1	L	125[A]	CYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	153	ASN	2.5
1	A	25	SER	2.5
1	D	151	MET	2.5
1	H	125[A]	CYS	2.4
1	E	25	SER	2.4
1	A	21	HIS	2.4
1	K	25	SER	2.4
1	D	1	MET	2.4
1	D	25	SER	2.3
1	C	41	ALA	2.3
1	I	145	LYS	2.3
1	L	2	LYS	2.3
1	E	37	GLN	2.3
1	F	18	ARG	2.3
1	G	145	LYS	2.3
1	F	151	MET	2.3
1	J	145	LYS	2.3
1	I	19	GLU	2.3
1	G	29	SER	2.3
1	K	22	ILE	2.3
1	J	23	TYR	2.2
1	F	25	SER	2.2
1	C	153	ASN	2.2
1	C	25	SER	2.2
1	I	45	GLU	2.1
1	H	18	ARG	2.1
1	J	18	ARG	2.1
1	B	125[A]	CYS	2.1
1	C	40	GLN	2.1
1	I	26	GLN	2.0
1	H	25	SER	2.0
1	K	125[A]	CYS	2.0
1	J	2	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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(Å ²)	Q<0.9
2	SO4	F	1155	5/5	0.76	0.13	41,55,67,68	0
2	SO4	H	1157	5/5	0.78	0.13	50,60,67,70	0
2	SO4	B	1148	5/5	0.81	0.10	57,58,71,74	0
2	SO4	D	1154	5/5	0.85	0.12	53,56,60,66	0
2	SO4	D	1153	5/5	0.86	0.11	43,52,63,67	0
2	SO4	F	1154	5/5	0.87	0.19	31,33,41,49	0
2	SO4	E	1148	5/5	0.90	0.14	33,47,55,56	0
2	SO4	C	1156	5/5	0.90	0.10	35,44,55,57	0
2	SO4	J	1149	5/5	0.90	0.25	39,40,46,52	0
2	SO4	J	1148	5/5	0.91	0.08	46,53,65,66	0
2	SO4	C	1155	5/5	0.92	0.14	31,36,49,53	0
2	SO4	K	1153	5/5	0.92	0.19	41,42,46,46	0
2	SO4	E	1149	5/5	0.93	0.16	39,42,49,50	0
2	SO4	H	1156	5/5	0.95	0.14	33,34,54,55	0
3	TRS	B	1149	8/8	0.95	0.07	14,15,16,17	0
3	TRS	K	1155	8/8	0.95	0.07	16,18,19,22	0
3	TRS	H	1158	8/8	0.96	0.06	15,16,18,19	0
3	TRS	E	1150	8/8	0.96	0.07	14,16,19,19	0
2	SO4	A	1148	5/5	0.97	0.08	21,29,36,38	0
2	SO4	L	1148	5/5	0.97	0.07	25,27,38,46	0
2	SO4	D	1152	5/5	0.97	0.08	17,27,33,43	0
2	SO4	J	1147	5/5	0.97	0.09	22,27,31,32	0
2	SO4	G	1147	5/5	0.97	0.09	23,27,39,44	0
2	SO4	H	1155	5/5	0.97	0.08	19,25,38,40	0
2	SO4	I	1147	5/5	0.98	0.06	22,29,38,39	0
2	SO4	K	1152	5/5	0.98	0.07	24,25,30,33	0
2	SO4	B	1147	5/5	0.98	0.07	18,21,36,38	0
2	SO4	F	1153	5/5	0.98	0.06	17,25,35,47	0
2	SO4	L	1147	5/5	0.99	0.04	17,18,20,25	0
2	SO4	F	1152	5/5	0.99	0.03	15,17,19,22	0
2	SO4	C	1154	5/5	0.99	0.07	18,23,29,33	0
2	SO4	H	1154	5/5	0.99	0.03	15,16,18,23	0
2	SO4	E	1147	5/5	0.99	0.05	20,24,28,29	0
2	SO4	K	1154	5/5	0.99	0.06	28,29,37,47	0
2	SO4	A	1147	5/5	1.00	0.03	13,13,16,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.