



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:49 AM UTC

PDB ID : 2IVQ / pdb_00002ivq
Title : SITE DIRECTED MUTAGENESIS OF KEY RESIDUES INVOLVED IN
THE CATALYTIC MECHANISM OF CYANASE
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Deposited on : 2006-06-14
Resolution : 2.10 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

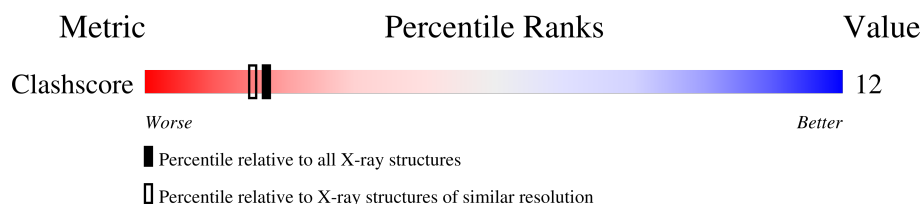
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
Clashscore	190562	7164 (2.10-2.10)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	156	82% 17% .
1	B	156	83% 16% .
1	C	156	78% 21% .
1	D	156	85% 15% .
1	E	156	80% 19% .
1	F	156	83% 16% .
1	G	156	78% 22% .
1	H	156	76% 22% .
1	I	156	83% 16% .
1	J	156	78% 21% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	F	1158	-	-	X	-
2	SO4	I	1158	-	-	X	-
3	CL	J	1160	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13229 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 CYANATE HYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	B	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	C	156	Total	C	N	O	S	0	1	0
			1202	772	198	227	5			
1	D	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	E	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	F	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	G	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	H	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	I	156	Total	C	N	O	S	0	0	0
			1196	767	197	227	5			
1	J	156	Total	C	N	O	S	0	1	0
			1202	772	198	227	5			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	96	LYS	ARG	engineered mutation	UNP P00816
B	96	LYS	ARG	engineered mutation	UNP P00816
C	96	LYS	ARG	engineered mutation	UNP P00816
D	96	LYS	ARG	engineered mutation	UNP P00816
E	96	LYS	ARG	engineered mutation	UNP P00816
F	96	LYS	ARG	engineered mutation	UNP P00816
G	96	LYS	ARG	engineered mutation	UNP P00816
H	96	LYS	ARG	engineered mutation	UNP P00816
I	96	LYS	ARG	engineered mutation	UNP P00816

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Chain	Residue	Modelled	Actual	Comment	Reference
J	96	LYS	ARG	engineered mutation	UNP P00816

- 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	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	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	G	1	Total	O	S	0	0
			5	4	1		
2	G	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	I	1	Total	O	S	0	0
			5	4	1		
2	I	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		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		
3	C	2	Total	Cl	0	0
			2	2		
3	D	1	Total	Cl	0	0
			1	1		
3	E	1	Total	Cl	0	0
			1	1		
3	F	2	Total	Cl	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total 1	Cl 1	0	0
3	H	1	Total 1	Cl 1	0	0
3	I	1	Total 1	Cl 1	0	0
3	J	2	Total 2	Cl 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	112	Total 112	O 112	0	0
4	B	130	Total 130	O 130	0	0
4	C	107	Total 107	O 107	0	0
4	D	114	Total 114	O 114	0	0
4	E	131	Total 131	O 131	0	0
4	F	116	Total 116	O 116	0	0
4	G	98	Total 98	O 98	0	0
4	H	108	Total 108	O 108	0	0
4	I	91	Total 91	O 91	0	0
4	J	117	Total 117	O 117	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

Note EDS was not executed.

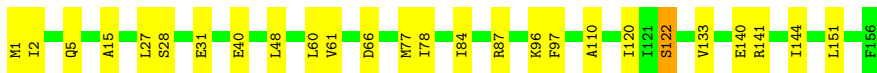
• Molecule 1: CYANATE HYDRATASE

Chain A:  82% 17%



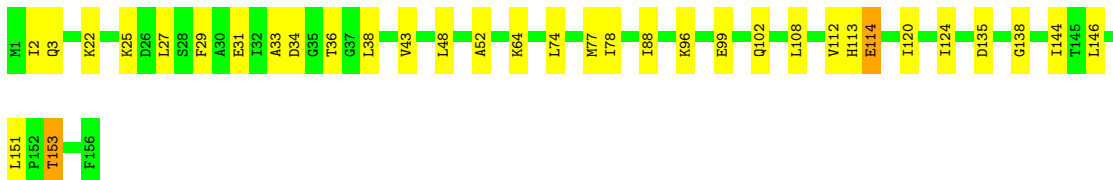
• Molecule 1: CYANATE HYDRATASE

Chain B:  83% 16%



• Molecule 1: CYANATE HYDRATASE

Chain C:  78% 21%



• Molecule 1: CYANATE HYDRATASE

Chain D:  85% 15%



• Molecule 1: CYANATE HYDRATASE

Chain E:  80% 19%



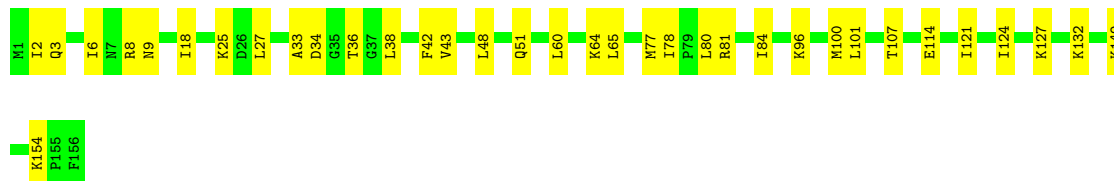
- Molecule 1: CYANATE HYDRATASE

Chain F:  83% 16%



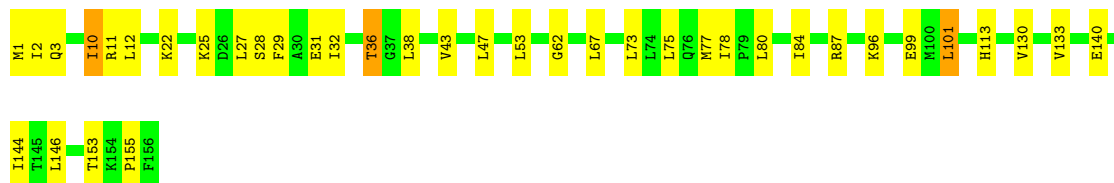
- Molecule 1: CYANATE HYDRATASE

Chain G:  78% 22%



- Molecule 1: CYANATE HYDRATASE

Chain H:  76% 22%



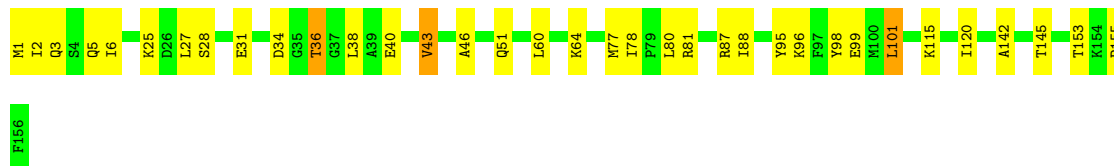
- Molecule 1: CYANATE HYDRATASE

Chain I:  83% 16%



- Molecule 1: CYANATE HYDRATASE

Chain J:  78% 21%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	76.88Å 80.84Å 82.65Å 69.99° 71.98° 66.17°	Depositor
Resolution (Å)	76.03 – 2.10	Depositor
% Data completeness (in resolution range)	95.7 (76.03-2.10)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.180 , 0.243	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13229	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CL

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	1.22	3/1213 (0.2%)	1.15	0/1638
1	B	1.27	1/1213 (0.1%)	1.10	2/1638 (0.1%)
1	C	1.17	0/1222	1.16	3/1649 (0.2%)
1	D	1.27	3/1213 (0.2%)	1.20	3/1638 (0.2%)
1	E	1.23	2/1213 (0.2%)	1.18	4/1638 (0.2%)
1	F	1.15	1/1213 (0.1%)	1.06	1/1638 (0.1%)
1	G	1.15	1/1213 (0.1%)	1.12	4/1638 (0.2%)
1	H	1.26	1/1213 (0.1%)	1.15	4/1638 (0.2%)
1	I	1.13	0/1213	1.07	2/1638 (0.1%)
1	J	1.29	3/1222 (0.2%)	1.13	2/1649 (0.1%)
All	All	1.22	15/12148 (0.1%)	1.13	25/16402 (0.2%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	6	ILE	N-CA	6.70	1.52	1.46
1	A	6	ILE	CA-CB	6.28	1.62	1.54
1	D	43	VAL	CA-CB	6.16	1.62	1.54
1	E	124	ILE	CA-CB	5.92	1.60	1.55
1	D	122	SER	CA-C	5.88	1.60	1.52
1	A	143	VAL	CA-CB	5.70	1.61	1.53
1	D	72	ILE	CA-CB	5.69	1.61	1.54
1	B	15	ALA	CA-CB	5.64	1.62	1.53
1	F	18	ILE	CA-CB	5.58	1.60	1.54
1	A	124	ILE	CA-CB	5.46	1.60	1.55
1	E	46	ALA	CA-CB	5.43	1.61	1.53
1	J	46	ALA	CA-CB	5.42	1.61	1.53
1	J	43	VAL	CA-CB	5.30	1.61	1.54
1	H	36	THR	CA-CB	5.28	1.61	1.54
1	G	100	MET	C-O	-5.26	1.17	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	135	ASP	CA-C-N	6.58	126.81	119.32
1	I	135	ASP	C-N-CA	6.58	126.81	119.32
1	E	10	ILE	N-CA-C	-6.50	104.42	110.53
1	H	10	ILE	N-CA-C	-6.47	104.45	110.53
1	J	36	THR	N-CA-C	-6.40	105.78	113.97
1	D	107	THR	N-CA-C	-6.35	104.44	111.36
1	H	36	THR	N-CA-C	-5.97	106.61	114.31
1	H	11	ARG	NE-CZ-NH2	5.85	124.46	119.20
1	D	60	LEU	N-CA-C	5.71	117.18	111.07
1	G	107	THR	N-CA-C	-5.52	105.34	111.36
1	J	101	LEU	N-CA-C	5.49	117.34	111.36
1	G	9	ASN	N-CA-C	5.42	117.94	111.71
1	C	153	THR	N-CA-C	-5.37	100.95	109.76
1	H	101	LEU	N-CA-C	5.37	117.22	111.36
1	C	114	GLU	CA-C-N	5.36	127.89	120.29
1	C	114	GLU	C-N-CA	5.36	127.89	120.29
1	E	124	ILE	N-CA-C	-5.35	106.45	111.48
1	G	8	ARG	N-CA-C	5.34	117.85	111.71
1	E	135	ASP	CA-C-N	5.33	126.51	119.84
1	E	135	ASP	C-N-CA	5.33	126.51	119.84
1	B	61	VAL	N-CA-C	-5.13	105.54	110.72
1	G	101	LEU	N-CA-C	5.10	116.92	111.36
1	B	122	SER	N-CA-C	5.07	117.65	110.50
1	D	122	SER	N-CA-C	5.01	117.50	110.23
1	F	128	LEU	N-CA-C	5.00	116.98	109.23

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	1196	0	1243	32	0
1	B	1196	0	1243	31	0
1	C	1202	0	1256	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1196	0	1243	23	0
1	E	1196	0	1243	37	0
1	F	1196	0	1243	29	0
1	G	1196	0	1243	39	0
1	H	1196	0	1243	50	0
1	I	1196	0	1243	35	0
1	J	1202	0	1256	49	0
2	A	10	0	0	0	0
2	B	10	0	0	1	0
2	C	5	0	0	0	0
2	D	15	0	0	0	0
2	E	10	0	0	0	0
2	F	15	0	0	2	0
2	G	15	0	0	1	0
2	H	15	0	0	0	0
2	I	15	0	0	2	0
2	J	10	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	2	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	2	0	0	0	0
3	G	1	0	0	1	0
3	H	1	0	0	0	0
3	I	1	0	0	1	0
3	J	2	0	0	2	0
4	A	112	0	0	4	0
4	B	130	0	0	8	0
4	C	107	0	0	5	0
4	D	114	0	0	2	0
4	E	131	0	0	5	0
4	F	116	0	0	4	0
4	G	98	0	0	3	0
4	H	108	0	0	4	0
4	I	91	0	0	4	0
4	J	117	0	0	4	0
All	All	13229	0	12456	307	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 12.

All (307) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1:MET:CE	1:D:2:ILE:HG22	1.62	1.29
1:J:1:MET:CE	1:J:2:ILE:HG22	1.62	1.29
1:A:1:MET:CE	1:A:2:ILE:HG22	1.64	1.26
1:J:2:ILE:HD11	1:J:78:ILE:HD11	1.29	1.13
1:D:1:MET:HE3	1:D:2:ILE:HG22	1.15	1.10
1:B:1:MET:CE	1:B:2:ILE:HG22	1.81	1.10
1:C:27:LEU:HD12	1:C:31:GLU:OE1	1.52	1.09
1:I:2:ILE:HD11	1:I:78:ILE:HD11	1.33	1.08
1:H:2:ILE:HD11	1:H:78:ILE:CD1	1.84	1.08
1:J:1:MET:HE2	1:J:2:ILE:HG22	1.25	1.07
1:H:2:ILE:HD11	1:H:78:ILE:HD11	1.11	1.07
1:B:1:MET:HE2	1:B:2:ILE:HG22	1.19	1.06
1:G:2:ILE:HD11	1:G:78:ILE:HD11	1.32	1.06
1:F:2:ILE:HD11	1:F:78:ILE:HD11	1.11	1.06
1:H:1:MET:HE3	1:H:2:ILE:HG22	1.28	1.05
1:B:60:LEU:HD13	4:B:2047:HOH:O	1.55	1.05
1:B:2:ILE:HD11	1:B:78:ILE:HD11	1.39	1.04
1:F:2:ILE:HD11	1:F:78:ILE:CD1	1.90	1.01
1:E:2:ILE:HD11	1:E:78:ILE:HD11	1.41	0.99
1:A:1:MET:HE2	1:A:2:ILE:HG22	1.46	0.97
1:C:2:ILE:HD11	1:C:78:ILE:HD11	1.45	0.97
1:E:114:GLU:HG3	1:F:1:MET:HE1	1.45	0.97
1:B:1:MET:HE3	1:B:2:ILE:H	1.36	0.90
1:J:96[A]:LYS:NZ	1:J:99:GLU:OE2	2.02	0.90
1:I:101:LEU:HD21	4:J:2020:HOH:O	1.72	0.89
1:J:1:MET:HE3	1:J:2:ILE:HG22	1.55	0.89
1:C:114:GLU:HG3	1:I:1:MET:CE	2.03	0.89
1:F:2:ILE:CD1	1:F:78:ILE:HD11	2.01	0.89
1:A:1:MET:CE	1:A:2:ILE:CG2	2.52	0.88
1:A:1:MET:HE3	1:A:2:ILE:HG22	1.55	0.87
4:A:2091:HOH:O	1:J:153:THR:HG23	1.75	0.85
1:A:1:MET:HE2	1:A:2:ILE:CG2	2.05	0.84
1:C:36:THR:HG22	1:C:38:LEU:H	1.44	0.83
1:G:2:ILE:HD11	1:G:78:ILE:CD1	2.09	0.82
1:B:1:MET:HE3	4:B:2004:HOH:O	1.80	0.82
1:B:1:MET:HE2	1:B:2:ILE:CG2	2.07	0.82
1:C:48:LEU:HD22	1:G:6:ILE:HD13	1.63	0.80
1:B:2:ILE:HD11	1:B:78:ILE:CD1	2.09	0.80
1:G:36:THR:HG21	1:G:43:VAL:HG21	1.64	0.79
1:D:6:ILE:HD13	1:F:48:LEU:HD22	1.63	0.79
1:J:1:MET:HE2	1:J:2:ILE:CG2	2.12	0.79
1:D:1:MET:HE2	1:D:2:ILE:HG22	1.63	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:114:GLU:HG3	1:F:1:MET:CE	2.16	0.76
1:I:96:LYS:HG3	1:J:120:ILE:HD12	1.66	0.76
1:H:1:MET:CE	1:H:2:ILE:HG22	2.15	0.75
1:F:1:MET:HE3	1:F:2:ILE:HG22	1.69	0.75
1:H:2:ILE:CD1	1:H:78:ILE:HD11	2.06	0.75
1:G:114:GLU:HG3	1:H:1:MET:CE	2.17	0.74
1:D:1:MET:HE3	1:D:2:ILE:CG2	2.08	0.73
1:H:36:THR:HG21	1:H:43:VAL:HG21	1.70	0.73
1:I:2:ILE:CD1	1:I:78:ILE:HD11	2.16	0.72
1:F:140:GLU:HG2	4:F:2105:HOH:O	1.88	0.72
1:J:36:THR:HG21	1:J:43:VAL:HG21	1.71	0.72
1:A:96:LYS:HG3	1:D:120:ILE:HD12	1.71	0.72
1:J:36:THR:HG22	1:J:38:LEU:H	1.55	0.72
1:G:2:ILE:HG21	1:H:113:HIS:O	1.90	0.71
1:A:100:MET:HB3	4:A:2080:HOH:O	1.91	0.70
1:E:36:THR:HG21	1:E:43:VAL:HG21	1.72	0.70
1:H:96:LYS:HE2	1:J:96[B]:LYS:HZ3	1.55	0.70
1:J:2:ILE:HD11	1:J:78:ILE:CD1	2.16	0.70
1:B:48:LEU:HD22	1:E:6:ILE:HD13	1.74	0.69
1:H:73:LEU:HD12	4:H:2059:HOH:O	1.93	0.69
1:A:6:ILE:HD13	1:I:48:LEU:HD22	1.74	0.68
1:G:36:THR:HG22	1:G:38:LEU:H	1.57	0.68
1:C:52:ALA:O	4:C:2040:HOH:O	2.10	0.68
1:C:124:ILE:HD12	4:I:2088:HOH:O	1.93	0.67
1:H:96:LYS:NZ	1:H:99:GLU:OE2	2.28	0.67
1:B:1:MET:HE3	1:B:2:ILE:N	2.07	0.66
1:J:36:THR:CG2	1:J:38:LEU:H	2.08	0.66
1:E:2:ILE:CD1	1:E:78:ILE:HD11	2.20	0.66
4:E:2060:HOH:O	1:G:127:LYS:HD2	1.95	0.65
1:E:120:ILE:HD12	1:G:96:LYS:HG3	1.77	0.65
1:C:124:ILE:CD1	4:I:2088:HOH:O	2.44	0.65
1:H:28:SER:HB2	1:I:1:MET:HE2	1.79	0.64
1:H:133:VAL:O	1:H:133:VAL:HG23	1.97	0.64
1:C:36:THR:CG2	1:C:38:LEU:HB2	2.27	0.64
3:C:1158:CL:CL	3:J:1160:CL:CL	2.89	0.64
1:G:114:GLU:HG3	1:H:1:MET:HE2	1.78	0.64
1:H:96:LYS:HZ1	1:J:96[B]:LYS:HE2	1.62	0.64
1:A:1:MET:HE3	1:A:2:ILE:H	1.63	0.64
1:C:114:GLU:HG3	1:I:1:MET:HE1	1.77	0.64
1:C:2:ILE:HD11	1:C:78:ILE:CD1	2.23	0.63
1:H:96:LYS:CE	1:J:96[B]:LYS:NZ	2.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:ILE:CD1	1:B:78:ILE:HD11	2.20	0.63
1:F:18:ILE:HG23	1:F:65:LEU:HD13	1.81	0.63
1:E:88:ILE:HD12	1:E:98:TYR:CZ	2.35	0.61
1:G:80:LEU:C	1:G:80:LEU:HD13	2.24	0.61
1:D:1:MET:CE	1:D:2:ILE:CG2	2.57	0.61
1:B:1:MET:CE	4:B:2004:HOH:O	2.41	0.61
1:E:36:THR:HG23	1:E:38:LEU:HB2	1.83	0.60
1:C:27:LEU:CD1	1:C:31:GLU:OE1	2.41	0.60
1:F:108:LEU:O	1:F:112:VAL:HG23	2.01	0.59
1:J:3:GLN:HB3	1:J:77:MET:HE3	1.84	0.59
1:G:33:ALA:O	1:G:36:THR:HB	2.02	0.59
1:H:96:LYS:HE2	1:J:96[B]:LYS:NZ	2.18	0.59
1:I:120:ILE:HD11	3:J:1160:CL:CL	2.40	0.58
1:G:124:ILE:HG23	1:H:153:THR:OG1	2.04	0.57
1:I:122:SER:HB2	1:I:151:LEU:HD11	1.85	0.57
1:A:1:MET:HE3	4:A:2001:HOH:O	2.03	0.57
1:G:51:GLN:NE2	1:G:81:ARG:HH21	2.02	0.57
1:C:113:HIS:HB3	1:I:78:ILE:HD12	1.86	0.57
1:E:36:THR:HG22	1:E:38:LEU:H	1.69	0.57
1:I:132:LYS:HE2	1:J:115:LYS:O	2.05	0.57
1:F:2:ILE:CD1	1:F:78:ILE:CD1	2.74	0.57
1:G:114:GLU:HG3	1:H:1:MET:HE1	1.86	0.57
1:C:114:GLU:HG3	1:I:1:MET:HE3	1.85	0.56
1:H:133:VAL:O	1:H:133:VAL:CG2	2.53	0.56
1:E:78:ILE:HD12	1:F:113:HIS:HB3	1.87	0.56
1:C:33:ALA:O	1:C:36:THR:HB	2.05	0.56
1:J:2:ILE:CD1	1:J:78:ILE:HD11	2.19	0.56
2:I:1158:SO4:O3	1:J:87:ARG:HD2	2.05	0.56
1:E:128:LEU:C	1:E:128:LEU:HD23	2.31	0.55
1:I:101:LEU:CD2	4:J:2020:HOH:O	2.42	0.55
1:J:36:THR:HG23	1:J:38:LEU:HB2	1.87	0.55
1:A:51:GLN:NE2	1:A:81:ARG:HH21	2.04	0.55
1:B:40:GLU:N	2:B:1157:SO4:O2	2.35	0.55
4:C:2032:HOH:O	1:H:101:LEU:HB3	2.06	0.55
1:D:51:GLN:NE2	1:D:81:ARG:HH21	2.04	0.55
1:G:36:THR:HG21	1:G:43:VAL:CG2	2.35	0.55
1:A:146:LEU:HD11	1:D:146:LEU:HD11	1.89	0.55
1:C:25:LYS:CB	1:C:27:LEU:HD22	2.37	0.55
1:C:96[A]:LYS:NZ	1:C:99:GLU:OE2	2.30	0.54
1:C:113:HIS:O	1:I:2:ILE:HG21	2.08	0.54
1:H:28:SER:CB	1:I:1:MET:HE2	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:ILE:HG13	1:C:151:LEU:HD12	1.88	0.54
1:B:27:LEU:HA	1:B:31:GLU:OE2	2.08	0.54
1:A:1:MET:HE1	1:A:2:ILE:HG22	1.80	0.54
1:D:56:ASP:OD1	1:D:59:ARG:NH2	2.41	0.53
1:B:1:MET:CE	1:B:2:ILE:H	2.15	0.53
1:J:3:GLN:CB	1:J:77:MET:HE3	2.38	0.53
1:E:88:ILE:HG13	4:G:2059:HOH:O	2.08	0.53
1:H:96:LYS:HE3	1:J:96[B]:LYS:HZ1	1.74	0.53
1:G:60:LEU:HD22	4:G:2030:HOH:O	2.08	0.53
1:C:34:ASP:O	1:C:64:LYS:NZ	2.42	0.53
1:H:36:THR:HG23	1:H:38:LEU:HB2	1.91	0.52
1:F:6:ILE:HD13	1:G:48:LEU:HD22	1.89	0.52
1:J:80:LEU:C	1:J:80:LEU:HD13	2.35	0.52
1:D:154:LYS:HB2	4:D:2114:HOH:O	2.08	0.52
1:C:25:LYS:HB2	1:C:27:LEU:HD22	1.91	0.52
1:C:146:LEU:HD11	1:H:146:LEU:HD11	1.92	0.52
1:H:2:ILE:CD1	1:H:78:ILE:CD1	2.74	0.52
1:A:142:ALA:CB	1:D:121:ILE:HD12	2.40	0.52
1:H:96:LYS:NZ	1:J:96[B]:LYS:HE2	2.24	0.52
1:A:31:GLU:O	1:A:34:ASP:HB2	2.10	0.51
1:E:36:THR:CG2	1:E:38:LEU:H	2.23	0.51
1:F:51:GLN:NE2	1:F:81:ARG:HH21	2.08	0.51
1:E:124:ILE:CD1	4:E:2044:HOH:O	2.58	0.51
1:B:28:SER:HB3	1:E:1:MET:HE3	1.93	0.51
1:G:38:LEU:HD22	1:H:155:PRO:HB3	1.91	0.51
1:G:2:ILE:CD1	1:G:78:ILE:CD1	2.88	0.51
1:B:96:LYS:HG3	1:F:120:ILE:HD12	1.93	0.51
1:G:36:THR:CG2	1:G:38:LEU:H	2.22	0.51
1:H:28:SER:OG	1:H:31:GLU:HG3	2.10	0.50
3:I:1160:CL:CL	1:J:96[A]:LYS:HE2	2.48	0.50
1:B:133:VAL:C	4:B:2119:HOH:O	2.53	0.50
1:E:135:ASP:OD1	1:G:149:LYS:CD	2.59	0.50
1:F:1:MET:HE3	1:F:2:ILE:CG2	2.39	0.50
1:H:10:ILE:HD13	4:H:2025:HOH:O	2.12	0.50
1:B:1:MET:CE	1:B:2:ILE:N	2.73	0.50
1:B:97:PHE:HE2	1:B:144:ILE:HD12	1.76	0.50
4:B:2105:HOH:O	1:D:1:MET:HE1	2.12	0.50
1:F:1:MET:CE	1:F:2:ILE:HG22	2.39	0.50
1:C:96[B]:LYS:HE2	4:E:2107:HOH:O	2.12	0.49
1:F:34:ASP:HB2	4:F:2037:HOH:O	2.12	0.49
1:C:135:ASP:HB3	1:C:138:GLY:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:34:ASP:O	1:J:64:LYS:NZ	2.45	0.49
1:A:38:LEU:HD22	1:J:155:PRO:CB	2.42	0.49
1:C:108:LEU:O	1:C:112:VAL:HG23	2.11	0.49
1:D:152:PRO:HG2	1:D:154:LYS:HE2	1.94	0.49
1:B:122:SER:HB2	1:B:151:LEU:HD11	1.93	0.49
1:C:2:ILE:HG22	4:C:2002:HOH:O	2.11	0.49
1:H:22:LYS:NZ	1:I:3:GLN:OE1	2.46	0.49
1:A:14:LEU:HD11	1:A:71:SER:HB3	1.94	0.49
4:A:2079:HOH:O	1:F:96:LYS:HG2	2.13	0.49
1:E:97:PHE:HE2	1:E:144:ILE:HD13	1.78	0.49
1:H:53:LEU:HD12	1:H:75:LEU:HD13	1.95	0.49
1:C:36:THR:HG21	1:C:43:VAL:HG21	1.94	0.48
1:D:25:LYS:CB	1:D:27:LEU:HD13	2.43	0.48
1:A:1:MET:HE2	1:A:2:ILE:HG23	1.93	0.48
1:A:35:GLY:O	1:A:60:LEU:HD22	2.13	0.48
1:A:128:LEU:C	1:A:128:LEU:HD12	2.37	0.48
1:G:2:ILE:CD1	1:G:78:ILE:HD11	2.23	0.48
1:E:77:MET:HG2	4:E:2086:HOH:O	2.13	0.48
1:A:38:LEU:HD22	1:J:155:PRO:HB3	1.94	0.48
1:B:141:ARG:HD3	4:B:2120:HOH:O	2.13	0.48
1:E:87:ARG:HD3	2:G:1158:SO4:O1	2.13	0.48
1:E:102:GLN:HE22	1:G:84:ILE:H	1.61	0.48
1:F:4:SER:N	1:F:77:MET:HE3	2.29	0.48
1:I:123:ALA:HB3	1:J:96[A]:LYS:HE3	1.96	0.48
1:F:149:LYS:NZ	4:F:2112:HOH:O	2.46	0.47
1:D:77:MET:HG2	4:D:2057:HOH:O	2.15	0.47
1:E:36:THR:CG2	1:E:38:LEU:HB2	2.45	0.47
1:E:128:LEU:HD23	1:E:129:ASP:N	2.29	0.47
1:G:42:PHE:CE1	1:H:155:PRO:HA	2.50	0.47
1:A:1:MET:HE3	1:A:2:ILE:N	2.27	0.47
1:C:2:ILE:CD1	1:C:78:ILE:HD11	2.31	0.47
4:C:2032:HOH:O	1:H:101:LEU:CD1	2.63	0.47
1:E:122:SER:HB2	1:E:151:LEU:HD11	1.97	0.47
1:H:140:GLU:HG2	4:H:2099:HOH:O	2.15	0.47
1:J:60:LEU:HD22	4:J:2043:HOH:O	2.14	0.47
1:C:74:LEU:HA	1:C:77:MET:HG3	1.97	0.47
1:G:36:THR:CG2	1:G:38:LEU:HB2	2.45	0.46
1:D:5:GLN:HG2	1:D:77:MET:HE2	1.96	0.46
1:D:36:THR:HG22	1:D:60:LEU:HD23	1.96	0.46
1:G:77:MET:HE2	1:G:77:MET:HB3	1.88	0.46
1:C:22:LYS:HE3	1:C:29:PHE:CE1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:80:LEU:C	1:G:80:LEU:CD1	2.87	0.46
1:H:36:THR:HG22	1:H:38:LEU:H	1.80	0.46
1:C:36:THR:CG2	1:C:38:LEU:H	2.21	0.46
1:C:114:GLU:HA	1:I:2:ILE:CG2	2.45	0.46
1:C:153:THR:OG1	1:I:124:ILE:HG23	2.14	0.46
1:E:135:ASP:OD1	1:G:149:LYS:HD3	2.16	0.46
2:I:1158:SO4:O3	1:J:87:ARG:CD	2.63	0.46
1:J:51:GLN:NE2	1:J:81:ARG:HH21	2.13	0.46
1:B:5:GLN:OE1	1:B:77:MET:HE2	2.16	0.46
1:E:88:ILE:HD12	1:E:98:TYR:CE1	2.51	0.45
1:E:142:ALA:HB2	1:G:121:ILE:HD12	1.98	0.45
1:I:122:SER:HG	1:I:125:ASN:N	2.13	0.45
1:C:96[B]:LYS:NZ	1:E:96:LYS:NZ	2.65	0.45
1:J:25:LYS:CB	1:J:27:LEU:HD13	2.46	0.45
1:C:22:LYS:HE3	1:C:29:PHE:CD1	2.51	0.45
1:G:3:GLN:HB3	1:G:77:MET:HE3	1.98	0.45
1:G:18:ILE:HG23	1:G:65:LEU:HD13	1.98	0.45
1:J:36:THR:CG2	1:J:38:LEU:HB2	2.46	0.45
1:H:32:ILE:HG21	1:H:47:LEU:HD11	1.99	0.45
1:A:102:GLN:NE2	1:D:84:ILE:H	2.15	0.44
1:A:120:ILE:HD11	1:A:151:LEU:HD12	2.00	0.44
1:B:120:ILE:HD12	1:F:96:LYS:HG3	1.98	0.44
1:C:96[B]:LYS:HZ3	1:E:96:LYS:HZ2	1.65	0.44
1:H:25:LYS:O	1:H:27:LEU:HD13	2.16	0.44
1:I:88:ILE:HD11	1:J:87:ARG:HB3	2.00	0.44
1:G:25:LYS:HB3	1:G:27:LEU:HD13	1.99	0.44
1:H:25:LYS:O	1:H:27:LEU:CD1	2.66	0.44
1:I:22:LYS:HE3	1:I:29:PHE:CE1	2.53	0.44
1:J:1:MET:HE3	1:J:2:ILE:N	2.32	0.44
1:J:40:GLU:N	2:J:1157:SO4:O3	2.46	0.44
1:C:96[B]:LYS:NZ	1:E:96:LYS:HD3	2.32	0.44
1:I:122:SER:OG	1:I:125:ASN:N	2.50	0.44
1:C:2:ILE:HD13	1:I:113:HIS:O	2.18	0.44
1:H:25:LYS:C	1:H:27:LEU:HD13	2.43	0.44
1:H:80:LEU:C	1:H:80:LEU:HD13	2.43	0.44
1:E:51:GLN:NE2	1:E:81:ARG:HH21	2.15	0.43
1:F:31:GLU:O	1:F:64:LYS:NZ	2.48	0.43
1:C:36:THR:HG23	1:C:38:LEU:HB2	2.00	0.43
1:A:135:ASP:O	1:A:138:GLY:N	2.43	0.43
1:F:3:GLN:C	1:F:77:MET:HE3	2.43	0.43
3:C:1159:CL:CL	3:G:1160:CL:CL	3.11	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:96:LYS:HE3	1:J:96[B]:LYS:NZ	2.30	0.43
1:J:1:MET:HE3	1:J:2:ILE:H	1.84	0.43
1:E:80:LEU:C	1:E:80:LEU:HD13	2.43	0.43
1:F:34:ASP:CB	4:F:2037:HOH:O	2.67	0.43
1:I:51:GLN:NE2	1:I:81:ARG:HH21	2.17	0.43
1:I:143:VAL:HG13	1:J:145:THR:CG2	2.49	0.43
1:C:77:MET:HG2	4:C:2056:HOH:O	2.19	0.43
1:C:78:ILE:HD12	1:I:113:HIS:HB3	2.01	0.43
1:A:142:ALA:HB2	1:D:121:ILE:HD12	2.00	0.43
1:A:4:SER:C	1:A:77:MET:HE2	2.44	0.42
1:B:97:PHE:CE2	1:B:144:ILE:HD12	2.54	0.42
1:G:77:MET:HG2	4:G:2044:HOH:O	2.19	0.42
1:B:77:MET:HE3	1:B:77:MET:HB3	1.88	0.42
1:H:3:GLN:HB3	1:H:77:MET:HE3	2.01	0.42
1:H:22:LYS:HE3	1:H:29:PHE:CE1	2.54	0.42
1:D:154:LYS:HA	1:D:155:PRO:HD3	1.93	0.42
1:B:110:ALA:HB1	1:F:29:PHE:CE1	2.54	0.42
1:C:144:ILE:HD12	1:C:144:ILE:N	2.34	0.42
1:I:121:ILE:HD12	1:J:142:ALA:CB	2.50	0.42
1:I:135:ASP:OD1	4:I:2083:HOH:O	2.22	0.42
1:D:1:MET:HE2	1:D:2:ILE:CG2	2.38	0.42
1:G:36:THR:HG23	1:G:38:LEU:HB2	2.02	0.42
1:E:33:ALA:O	1:E:36:THR:HB	2.20	0.42
1:F:87:ARG:NH1	2:F:1158:SO4:O4	2.53	0.42
1:A:80:LEU:HD13	1:A:80:LEU:C	2.45	0.41
1:A:142:ALA:HB3	1:D:121:ILE:HD12	2.01	0.41
1:C:25:LYS:HB2	1:C:27:LEU:CD2	2.49	0.41
1:A:42:PHE:CE1	1:J:155:PRO:HA	2.55	0.41
1:B:84:ILE:H	1:F:102:GLN:HE22	1.67	0.41
1:C:25:LYS:HB3	1:C:27:LEU:HD22	2.02	0.41
1:E:102:GLN:NE2	1:G:84:ILE:H	2.18	0.41
1:H:12:LEU:HD23	1:H:12:LEU:HA	1.93	0.41
1:B:87:ARG:HD2	2:F:1158:SO4:O2	2.20	0.41
1:H:62:GLY:HA2	1:H:67:LEU:HD12	2.02	0.41
1:I:130:VAL:HG22	1:I:144:ILE:HG12	2.02	0.41
1:A:77:MET:HE2	1:A:77:MET:HB3	1.91	0.41
1:J:5:GLN:NE2	4:J:2010:HOH:O	2.52	0.41
1:J:28:SER:OG	1:J:31:GLU:HG3	2.20	0.41
1:E:124:ILE:HD13	4:E:2044:HOH:O	2.21	0.41
1:C:77:MET:HB3	1:C:77:MET:HE2	1.76	0.41
1:H:130:VAL:HG22	1:H:144:ILE:HG12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:88:ILE:HD12	1:J:98:TYR:CZ	2.56	0.41
1:G:34:ASP:O	1:G:64:LYS:NZ	2.54	0.41
1:C:88:ILE:HD11	1:H:87:ARG:HB3	2.02	0.41
1:C:102:GLN:NE2	1:H:84:ILE:H	2.19	0.41
1:E:116:PHE:CE1	1:G:132:LYS:HG3	2.56	0.41
1:G:154:LYS:HD2	4:H:2042:HOH:O	2.20	0.41
1:I:121:ILE:HD12	1:J:142:ALA:HB2	2.02	0.41
1:I:97:PHE:HB3	1:J:101:LEU:HD21	2.03	0.41
1:A:80:LEU:HD22	1:A:80:LEU:HA	1.91	0.40
1:B:140:GLU:HG2	4:B:2119:HOH:O	2.20	0.40
1:C:102:GLN:HE22	1:H:84:ILE:H	1.69	0.40
1:J:95:TYR:CD2	1:J:95:TYR:C	2.99	0.40
1:C:114:GLU:HA	1:I:2:ILE:HG21	2.03	0.40
1:F:77:MET:HE2	1:F:77:MET:HB3	1.81	0.40
1:B:66:ASP:HB2	4:B:2043:HOH:O	2.22	0.40
1:C:2:ILE:HD13	1:C:2:ILE:HG21	1.85	0.40
1:C:3:GLN:HB3	1:C:77:MET:HE3	2.04	0.40
1:E:78:ILE:HD13	1:E:78:ILE:HG21	1.80	0.40
1:I:51:GLN:NE2	4:I:2042:HOH:O	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 37 ligands modelled in this entry, 13 are monoatomic - leaving 24 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	$\# Z > 2$
2	SO4	J	1157	-	4,4,4	0.53	0	6,6,6	0.50	0
2	SO4	J	1158	-	4,4,4	0.30	0	6,6,6	0.77	0
2	SO4	F	1159	-	4,4,4	0.33	0	6,6,6	0.28	0
2	SO4	D	1159	-	4,4,4	0.31	0	6,6,6	0.34	0
2	SO4	I	1158	-	4,4,4	0.31	0	6,6,6	1.87	2 (33%)
2	SO4	G	1159	-	4,4,4	0.26	0	6,6,6	0.16	0
2	SO4	I	1157	-	4,4,4	0.23	0	6,6,6	0.33	0
2	SO4	F	1157	-	4,4,4	0.23	0	6,6,6	0.38	0
2	SO4	A	1157	-	4,4,4	0.21	0	6,6,6	1.15	0
2	SO4	F	1158	-	4,4,4	0.59	0	6,6,6	1.43	1 (16%)
2	SO4	H	1159	-	4,4,4	0.24	0	6,6,6	0.53	0
2	SO4	B	1157	-	4,4,4	0.23	0	6,6,6	0.39	0
2	SO4	A	1158	-	4,4,4	0.25	0	6,6,6	0.22	0
2	SO4	D	1157	-	4,4,4	0.37	0	6,6,6	0.54	0
2	SO4	B	1158	-	4,4,4	0.23	0	6,6,6	0.15	0
2	SO4	E	1157	-	4,4,4	0.26	0	6,6,6	0.51	0
2	SO4	H	1157	-	4,4,4	0.32	0	6,6,6	0.45	0
2	SO4	C	1157	-	4,4,4	0.33	0	6,6,6	0.52	0
2	SO4	E	1158	-	4,4,4	0.29	0	6,6,6	0.39	0
2	SO4	G	1158	-	4,4,4	0.63	0	6,6,6	1.66	1 (16%)
2	SO4	H	1158	-	4,4,4	0.41	0	6,6,6	1.37	1 (16%)
2	SO4	D	1158	-	4,4,4	0.39	0	6,6,6	1.04	0
2	SO4	I	1159	-	4,4,4	0.36	0	6,6,6	0.33	0
2	SO4	G	1157	-	4,4,4	0.30	0	6,6,6	0.55	0

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	1158	SO4	O3-S-O1	-3.17	92.97	109.56
2	H	1158	SO4	O3-S-O2	-2.97	94.02	109.56
2	F	1158	SO4	O4-S-O2	-2.55	96.23	109.56
2	I	1158	SO4	O4-S-O1	2.34	121.81	109.56
2	G	1158	SO4	O4-S-O3	2.17	120.52	108.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	1157	SO4	1	0
2	I	1158	SO4	2	0
2	F	1158	SO4	2	0
2	B	1157	SO4	1	0
2	G	1158	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.