



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:28 PM UTC

PDB ID : 4MWA / pdb_00004mwa
Title : 1.85 Angstrom Crystal Structure of GCPE Protein from Bacillus anthracis
Authors : Minasov, G.; Wawrzak, Z.; Brunzelle, J.S.; Xu, X.; Cui, H.; Maltseva, N.; Bishop, B.; Kwon, K.; Savchenko, A.; Joachimiak, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2013-09-24
Resolution : 1.85 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

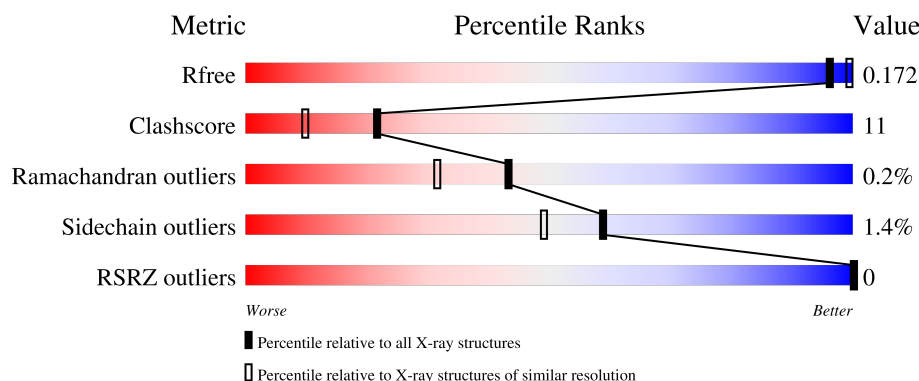
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	276	 69% 24% • 5%
1	G	276	 76% 16% • 7%
2	B	276	 79% 14% • 6%
3	C	276	 67% 25% • 7%
4	D	276	 67% 23% • 9%

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Mol	Chain	Length	Quality of chain
4	F	276	 72% 21% • 5%
4	H	276	 70% 22% • 8%
5	E	276	 78% 15% • 6%

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
6	SO4	E	401	-	-	X	-
6	SO4	G	402	-	-	X	-
6	SO4	H	401	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 17202 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 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	Se	0	10	0
			2083	1305	376	395	1	6			
1	G	258	Total	C	N	O	S	Se	0	0	0
			1965	1235	354	371	1	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q81LV7
A	-1	ASN	-	expression tag	UNP Q81LV7
A	0	ALA	-	expression tag	UNP Q81LV7
A	1	MSE	-	expression tag	UNP Q81LV7
A	2	ASN	-	expression tag	UNP Q81LV7
A	3	GLU	-	expression tag	UNP Q81LV7
G	-2	SER	-	expression tag	UNP Q81LV7
G	-1	ASN	-	expression tag	UNP Q81LV7
G	0	ALA	-	expression tag	UNP Q81LV7
G	1	MSE	-	expression tag	UNP Q81LV7
G	2	ASN	-	expression tag	UNP Q81LV7
G	3	GLU	-	expression tag	UNP Q81LV7

- Molecule 2 is a protein called 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	259	Total	C	N	O	S	Se	0	2	0
			1990	1250	360	375	1	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	SER	-	expression tag	UNP Q81LV7
B	-1	ASN	-	expression tag	UNP Q81LV7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	ALA	-	expression tag	UNP Q81LV7
B	1	MSE	-	expression tag	UNP Q81LV7
B	2	ASN	-	expression tag	UNP Q81LV7
B	3	GLU	-	expression tag	UNP Q81LV7

- Molecule 3 is a protein called 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	257	Total	C	N	O	S	Se	0	1	0
			1968	1236	354	373	1	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	SER	-	expression tag	UNP Q81LV7
C	-1	ASN	-	expression tag	UNP Q81LV7
C	0	ALA	-	expression tag	UNP Q81LV7
C	1	MSE	-	expression tag	UNP Q81LV7
C	2	ASN	-	expression tag	UNP Q81LV7
C	3	GLU	-	expression tag	UNP Q81LV7

- Molecule 4 is a protein called 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	251	Total	C	N	O	S	Se	0	4	0
			1938	1222	344	367	1	4			
4	F	262	Total	C	N	O	S	Se	0	5	0
			2033	1274	371	383	1	4			
4	H	255	Total	C	N	O	S	Se	0	1	0
			1951	1226	352	368	1	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	SER	-	expression tag	UNP Q81LV7
D	-1	ASN	-	expression tag	UNP Q81LV7
D	0	ALA	-	expression tag	UNP Q81LV7
D	1	MSE	-	expression tag	UNP Q81LV7
D	2	ASN	-	expression tag	UNP Q81LV7
D	3	GLU	-	expression tag	UNP Q81LV7
F	-2	SER	-	expression tag	UNP Q81LV7
F	-1	ASN	-	expression tag	UNP Q81LV7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	expression tag	UNP Q81LV7
F	1	MSE	-	expression tag	UNP Q81LV7
F	2	ASN	-	expression tag	UNP Q81LV7
F	3	GLU	-	expression tag	UNP Q81LV7
H	-2	SER	-	expression tag	UNP Q81LV7
H	-1	ASN	-	expression tag	UNP Q81LV7
H	0	ALA	-	expression tag	UNP Q81LV7
H	1	MSE	-	expression tag	UNP Q81LV7
H	2	ASN	-	expression tag	UNP Q81LV7
H	3	GLU	-	expression tag	UNP Q81LV7

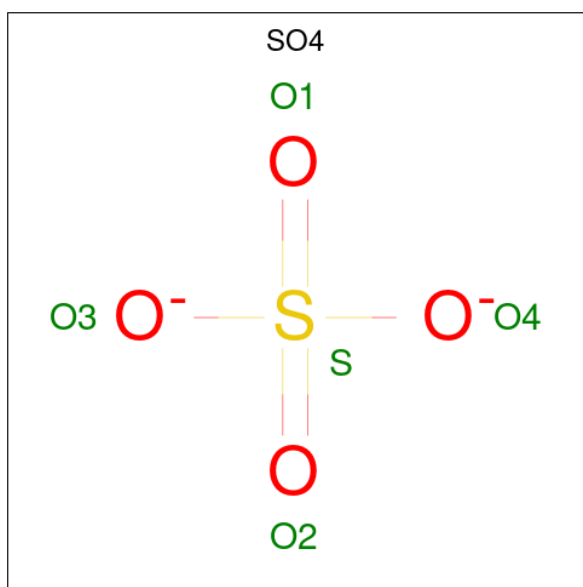
- Molecule 5 is a protein called 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
5	E	260	Total	C	N	O	S	Se	0	2	0
			2003	1256	363	379	1	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	SER	-	expression tag	UNP Q81LV7
E	-1	ASN	-	expression tag	UNP Q81LV7
E	0	ALA	-	expression tag	UNP Q81LV7
E	1	MSE	-	expression tag	UNP Q81LV7
E	2	ASN	-	expression tag	UNP Q81LV7
E	3	GLU	-	expression tag	UNP Q81LV7

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



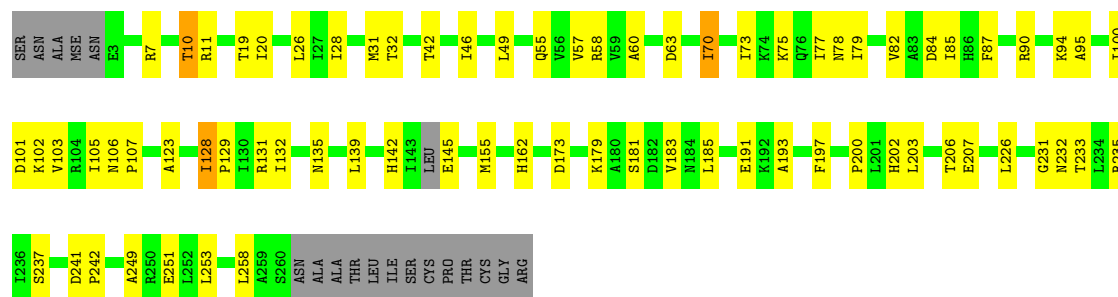
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	E	1	Total	O	S	0	0
			5	4	1		
6	E	1	Total	O	S	0	0
			5	4	1		
6	F	1	Total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

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

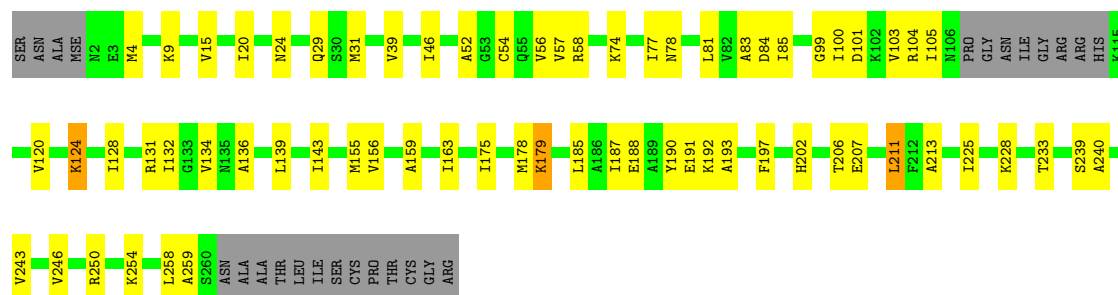
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	146	Total O 147 147	0	2
8	B	139	Total O 142 142	0	3
8	C	147	Total O 149 149	0	2
8	D	116	Total O 119 119	0	3
8	E	166	Total O 172 172	0	6
8	F	148	Total O 154 154	0	6
8	G	158	Total O 162 162	0	4
8	H	148	Total O 151 151	0	3



- Molecule 4: 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase

Chain D: 67% 23% 9%



- Molecule 4: 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase

Chain F: 72% 21% 5%



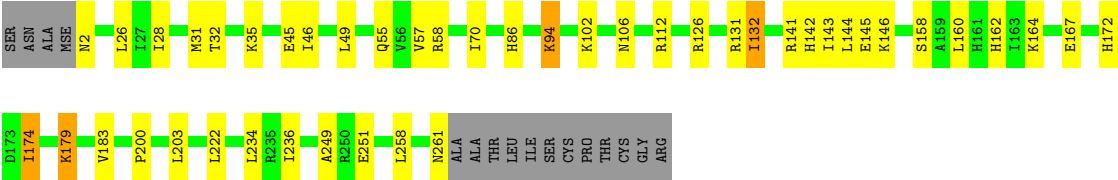
- Molecule 4: 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase

Chain H: 70% 22% 8%



- Molecule 5: 4-hydroxy-3-methylbut-2-en-1-yl diphosphate synthase

Chain E: 78% 15% • 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	162.68Å 162.68Å 76.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.77 – 1.85 29.77 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.4 (29.77-1.85) 99.4 (29.77-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.66 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.8.0046	Depositor
R, R_{free}	0.125 , 0.169 0.133 , 0.172	Depositor DCC
R_{free} test set	9677 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 17.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.468 for -h,-k,l 0.479 for h,-h-k,-l 0.468 for -k,-h,-l	Xtriage
Reported twinning fraction	0.241 for H, K, L 0.260 for -K, -H, -L 0.268 for -h,-k,l 0.232 for K, H, -L	Depositor
Outliers	0 of 192077 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	17202	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 3.77% 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: SO4, KCX, 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	0.71	2/2093 (0.1%)	0.98	3/2817 (0.1%)
1	G	0.76	0/1974	1.03	3/2660 (0.1%)
2	B	0.65	0/2000	0.97	3/2696 (0.1%)
3	C	0.66	0/1963	0.97	5/2643 (0.2%)
4	D	0.64	0/1957	0.97	3/2636 (0.1%)
4	F	0.71	0/2055	1.03	4/2770 (0.1%)
4	H	0.73	0/1973	1.02	7/2659 (0.3%)
5	E	0.74	0/2000	1.00	6/2694 (0.2%)
All	All	0.70	2/16015 (0.0%)	1.00	34/21575 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	48[A]	ARG	C-O	-5.80	1.17	1.24
1	A	48[B]	ARG	C-O	-5.80	1.17	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	70	ILE	N-CA-C	8.42	118.99	110.23
4	H	256	PHE	N-CA-C	8.21	123.34	113.16
1	A	110	ILE	N-CA-C	7.90	117.96	110.30
1	G	110	ILE	N-CA-C	7.82	118.59	110.62
4	D	39	VAL	N-CA-C	7.34	117.42	110.30
3	C	70	ILE	N-CA-C	6.88	117.39	110.23
3	C	183	VAL	N-CA-C	6.68	116.83	110.42
4	H	111	GLY	N-CA-C	6.61	119.58	111.45
5	E	174	ILE	N-CA-C	6.57	117.93	108.48
4	F	174	ILE	N-CA-C	6.53	117.88	108.48
4	H	106	ASN	CA-C-N	6.47	127.92	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	106	ASN	C-N-CA	6.47	127.92	119.84
1	A	174	ILE	N-CA-C	6.19	117.68	108.46
1	A	179	LYS	N-CA-C	5.99	119.28	109.76
4	F	179	LYS	N-CA-C	5.93	119.31	109.46
4	D	179	LYS	N-CA-C	5.87	119.10	109.76
1	G	183	VAL	N-CA-C	5.78	116.51	110.62
4	H	39	VAL	N-CA-C	5.69	115.88	110.42
5	E	86	HIS	N-CA-C	5.62	117.49	111.36
1	G	39	VAL	N-CA-C	5.49	115.69	110.42
5	E	179	LYS	N-CA-C	5.40	117.55	108.96
3	C	63	ASP	N-CA-C	5.37	115.22	108.45
4	H	128	ILE	N-CA-C	5.36	113.24	108.63
3	C	106	ASN	CA-C-N	5.33	126.51	119.84
3	C	106	ASN	C-N-CA	5.33	126.51	119.84
4	D	259	ALA	N-CA-C	5.33	116.80	110.19
5	E	132	ILE	CB-CA-C	-5.32	104.85	111.08
2	B	179	LYS	N-CA-C	5.29	117.38	108.96
4	F	196	ALA	N-CA-C	5.28	117.95	111.82
5	E	146	LYS	N-CA-C	5.28	117.11	111.36
4	F	35	LYS	N-CA-C	-5.26	102.07	109.96
2	B	74	LYS	N-CA-C	5.17	116.92	111.28
4	H	128	ILE	CB-CA-C	-5.13	106.17	110.94
2	B	76	GLN	N-CA-C	5.09	119.58	113.17

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	2083	0	2132	58	0
1	G	1965	0	2026	32	0
2	B	1990	0	2047	28	0
3	C	1968	0	2022	60	0
4	D	1938	0	2007	50	0
4	F	2033	0	2095	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1951	0	2011	52	0
5	E	2003	0	2054	32	0
6	A	5	0	0	1	0
6	B	5	0	0	1	0
6	C	10	0	0	0	0
6	D	10	0	0	0	0
6	E	10	0	0	2	0
6	F	10	0	0	1	0
6	G	10	0	0	2	0
6	H	5	0	0	2	0
7	A	1	0	0	1	0
7	B	3	0	0	0	0
7	C	3	0	0	0	0
7	F	3	0	0	1	0
8	A	147	0	0	4	0
8	B	142	0	0	4	0
8	C	149	0	0	4	0
8	D	119	0	0	2	0
8	E	172	0	0	7	0
8	F	154	0	0	3	0
8	G	162	0	0	3	0
8	H	151	0	0	4	0
All	All	17202	0	16394	351	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 11.

All (351) 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)
4:F:155:MSE:SE	4:F:185:LEU:HD21	2.10	1.02
4:H:236:ILE:HG22	4:H:238:LEU:CD1	1.95	0.97
4:H:29:GLN:HE22	4:H:58:ARG:NH1	1.68	0.91
3:C:241:ASP:HB3	8:C:554:HOH:O	1.73	0.87
4:H:28:ILE:HD11	4:H:253:LEU:HD12	1.61	0.82
5:E:28:ILE:HD13	5:E:234:LEU:HD21	1.62	0.81
4:F:250:ARG:HG3	4:F:250:ARG:HH11	1.45	0.80
4:D:46[A]:ILE:HD12	4:D:57:VAL:HG11	1.61	0.80
4:H:128:ILE:HD11	8:H:507:HOH:O	1.82	0.78
3:C:19:THR:C	3:C:20:ILE:HD12	2.08	0.78
4:F:5:THR:HG23	4:F:9:LYS:HG3	1.64	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:60:ALA:HA	3:C:84:ASP:HB3	1.66	0.77
4:H:177:SER:OG	4:H:202:HIS:HB3	1.84	0.77
4:H:236:ILE:HG22	4:H:238:LEU:HD11	1.65	0.77
3:C:70:ILE:HD12	3:C:100:ILE:HD11	1.65	0.76
4:D:187[B]:ILE:HA	4:D:225:ILE:HD11	1.68	0.75
1:A:140:GLU:HB2	1:A:143:ILE:HD13	1.69	0.75
3:C:10:THR:HG21	3:C:200:PRO:HA	1.67	0.74
2:B:141:ARG:NH1	8:B:586:HOH:O	2.21	0.74
4:D:187[A]:ILE:HA	4:D:225:ILE:HD11	1.73	0.71
3:C:46:ILE:HD12	3:C:57:VAL:HG11	1.73	0.70
4:D:31:MSE:SE	4:D:58:ARG:HH12	2.24	0.69
4:H:236:ILE:CG2	4:H:238:LEU:HD11	2.22	0.69
5:E:28:ILE:HD12	5:E:249:ALA:HB1	1.74	0.69
2:B:15:VAL:HG21	2:B:20:ILE:HD12	1.74	0.69
4:D:202:HIS:HA	4:D:233:THR:HG23	1.74	0.69
3:C:75:LYS:HG3	8:C:619:HOH:O	1.93	0.68
1:A:203[A]:LEU:HD11	1:A:231:GLY:HA3	1.75	0.68
1:A:241:ASP:OD2	1:A:243:VAL:HG12	1.94	0.68
1:A:228[A]:LYS:HE2	1:A:228[A]:LYS:HA	1.75	0.67
4:F:5:THR:HG23	4:F:9:LYS:CG	2.24	0.66
4:F:236:ILE:HG21	4:F:248:VAL:HG11	1.77	0.66
2:B:49:LEU:HD21	2:B:242:PRO:HB2	1.77	0.66
4:D:240:ALA:HB1	8:D:528:HOH:O	1.95	0.66
1:G:4:MSE:SE	1:G:228:LYS:HG2	2.46	0.66
4:D:120:VAL:O	4:D:124:LYS:HG2	1.96	0.65
1:G:211:LEU:HG	1:G:240:ALA:HB2	1.78	0.65
1:A:234:LEU:HD12	1:A:234:LEU:C	2.22	0.65
3:C:123:ALA:HA	3:C:128:ILE:HG23	1.79	0.65
5:E:45:GLU:O	5:E:49:LEU:HD13	1.98	0.64
1:G:49:LEU:HD21	1:G:242:PRO:HB2	1.78	0.64
5:E:160:LEU:O	5:E:164:LYS:HG3	1.97	0.64
1:G:88:ASP:CG	8:G:623:HOH:O	2.41	0.63
4:H:253:LEU:O	4:H:257:GLY:HA2	1.98	0.63
2:B:13:VAL:HG11	2:B:175:ILE:HD11	1.79	0.63
4:F:33:THR:HG22	4:F:45:GLU:OE2	1.98	0.62
5:E:179:LYS:HD2	6:E:401:SO4:O1	2.00	0.62
3:C:207:GLU:HG2	3:C:237:SER:HB3	1.82	0.62
4:F:185:LEU:O	4:F:185:LEU:HD23	2.00	0.62
4:D:31:MSE:SE	4:D:58:ARG:NH1	2.82	0.62
5:E:28:ILE:CD1	5:E:234:LEU:HD21	2.30	0.61
5:E:94[A]:LYS:HZ3	5:E:94[A]:LYS:HB3	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:89:TYR:HB3	2:B:110:ILE:HG23	1.82	0.61
3:C:20:ILE:HD12	3:C:20:ILE:N	2.15	0.61
1:A:159:ALA:O	1:A:163:ILE:HG12	2.02	0.60
4:F:254:LYS:NZ	8:F:611:HOH:O	2.33	0.60
4:H:20:ILE:CD1	4:H:27:ILE:HG13	2.31	0.60
1:A:250:ARG:O	1:A:254:LYS:HG3	2.02	0.59
3:C:90:ARG:O	3:C:94:LYS:HG2	2.02	0.59
1:A:106:ASN:HB2	1:A:109:ASN:ND2	2.18	0.59
1:A:188[B]:GLU:OE1	1:A:188[B]:GLU:HA	2.02	0.59
3:C:77:ILE:HG22	3:C:78:ASN:N	2.18	0.59
2:B:58:ARG:NH1	6:B:401:SO4:O2	2.35	0.59
5:E:32:THR:CG2	5:E:46:ILE:HD11	2.33	0.59
1:G:243:VAL:HG12	1:G:247:LYS:HE3	1.85	0.58
3:C:10:THR:HG21	3:C:200:PRO:CA	2.32	0.58
3:C:155:MSE:HE1	3:C:185:LEU:HD21	1.85	0.58
5:E:112:ARG:NH1	8:E:630[A]:HOH:O	2.36	0.58
4:D:155:MSE:HE1	4:D:185:LEU:HD21	1.86	0.58
4:H:58:ARG:CZ	4:H:104:ARG:HD3	2.34	0.57
3:C:103:VAL:HG23	3:C:128:ILE:HD11	1.84	0.57
4:H:56:VAL:HG22	4:H:80:PRO:HG2	1.86	0.57
4:D:178:MSE:HE3	4:D:193:ALA:HB2	1.86	0.57
4:H:107:PRO:O	4:H:110:ILE:HG22	2.04	0.57
4:F:250:ARG:HH11	4:F:250:ARG:CG	2.16	0.57
4:H:29:GLN:HG2	4:H:30:SER:N	2.20	0.57
1:G:4:MSE:SE	1:G:191:GLU:HG3	2.55	0.57
1:A:49:LEU:HD21	1:A:242:PRO:HB2	1.86	0.56
3:C:70:ILE:HG21	3:C:100:ILE:HG12	1.86	0.56
3:C:203:LEU:HD11	3:C:231:GLY:HA3	1.86	0.56
3:C:207:GLU:HG3	8:D:584:HOH:O	2.06	0.56
3:C:28:ILE:HD12	3:C:249:ALA:HB1	1.86	0.56
1:A:249:ALA:O	1:A:253:LEU:HD13	2.05	0.56
2:B:13:VAL:CG1	2:B:175:ILE:HD11	2.35	0.56
1:G:31:MSE:HE1	1:G:58:ARG:NH1	2.21	0.56
4:H:241:ASP:OD1	4:H:242:PRO:HD2	2.06	0.55
2:B:95:ALA:HB1	2:B:100:ILE:HG13	1.89	0.55
3:C:102:KCX:OQ2	3:C:131:ARG:NE	2.37	0.55
1:A:89[B]:TYR:HB2	1:A:119:VAL:CG2	2.36	0.55
1:G:102:KCX:OQ1	1:G:131:ARG:NE	2.35	0.55
4:D:211:LEU:HD23	4:D:211:LEU:C	2.31	0.55
5:E:55:GLN:NE2	5:E:261:ASN:OD1	2.40	0.55
4:F:201:LEU:HD22	4:F:230:ILE:CG2	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:58:ARG:NH1	6:H:401:SO4:O1	2.40	0.55
2:B:103:VAL:HG12	2:B:105:ILE:HG23	1.88	0.55
3:C:207:GLU:CG	3:C:237:SER:HB3	2.36	0.55
3:C:235:ARG:C	3:C:235:ARG:HD3	2.31	0.55
4:H:236:ILE:CG2	4:H:238:LEU:CD1	2.77	0.55
5:E:28:ILE:HD13	5:E:234:LEU:CD2	2.33	0.54
3:C:58:ARG:HG3	3:C:82:VAL:HB	1.89	0.54
3:C:155:MSE:CE	3:C:185:LEU:HD21	2.37	0.54
5:E:179:LYS:HG3	8:E:625:HOH:O	2.06	0.54
4:F:226:LEU:HD12	4:F:252:LEU:HD21	1.89	0.54
4:F:234:LEU:HD12	4:F:234:LEU:C	2.33	0.54
4:H:110:ILE:CG2	4:H:116:VAL:HG22	2.36	0.54
4:D:207:GLU:OE1	4:D:239:SER:OG	2.24	0.54
4:F:65:ARG:HG3	8:F:622:HOH:O	2.07	0.54
1:A:132:ILE:HD12	1:A:163:ILE:HD13	1.89	0.54
4:F:77:ILE:CD1	4:F:81:LEU:HD21	2.38	0.54
4:H:58:ARG:NH2	4:H:82:VAL:HG11	2.23	0.53
3:C:11:ARG:NH2	3:C:173:ASP:OD1	2.40	0.53
3:C:32:THR:CG2	3:C:46:ILE:HD11	2.39	0.53
1:A:90[B]:ARG:NH2	7:A:402:CL:CL	2.78	0.53
4:H:236:ILE:HG22	4:H:238:LEU:HD12	1.87	0.53
4:D:178:MSE:HE2	4:D:197:PHE:HZ	1.73	0.52
4:D:4:MSE:HE1	4:D:228:LYS:HB3	1.90	0.52
4:F:60:ALA:HA	4:F:84:ASP:HB3	1.91	0.52
4:H:64:GLU:HG3	4:H:94:LYS:HE2	1.91	0.52
3:C:135:ASN:ND2	8:C:510:HOH:O	2.28	0.52
1:G:4:MSE:SE	1:G:228:LYS:CG	3.07	0.52
1:G:7:ARG:NH1	1:G:226:LEU:O	2.41	0.52
1:A:26:LEU:HD11	1:A:258:LEU:HD13	1.91	0.52
4:H:11:ARG:NH2	4:H:173:ASP:OD1	2.42	0.52
4:H:96:ILE:HG23	4:H:128:ILE:CD1	2.39	0.52
4:H:179:LYS:NZ	8:H:506:HOH:O	2.42	0.52
3:C:49:LEU:HD21	3:C:242:PRO:HB2	1.90	0.52
4:F:185:LEU:HD23	4:F:185:LEU:C	2.34	0.52
1:G:27:ILE:HD11	8:G:539:HOH:O	2.09	0.52
4:D:77:ILE:HD11	4:D:81:LEU:HG	1.91	0.52
4:F:155:MSE:SE	4:F:185:LEU:CD2	2.99	0.52
1:A:25:GLU:N	1:A:25:GLU:OE1	2.43	0.52
4:F:251[B]:GLU:HG3	1:G:212:PHE:CE2	2.44	0.52
4:H:101:ASP:O	4:H:129:PRO:HD2	2.10	0.52
3:C:193:ALA:HB1	3:C:197:PHE:CZ	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:251:GLU:HG2	4:H:213:ALA:HA	1.91	0.51
4:H:29:GLN:HG2	4:H:30:SER:H	1.75	0.51
4:H:96:ILE:HG23	4:H:128:ILE:HD12	1.92	0.51
3:C:77:ILE:HG22	3:C:78:ASN:H	1.75	0.51
3:C:123:ALA:HA	3:C:128:ILE:CG2	2.40	0.51
4:F:152:ALA:O	4:F:156:VAL:HG23	2.11	0.51
1:A:89[A]:TYR:CE1	1:A:93:LEU:HD11	2.45	0.51
1:A:67:ALA:HB3	8:A:533:HOH:O	2.08	0.51
3:C:95:ALA:HB1	3:C:100:ILE:HG13	1.93	0.51
1:G:108:GLY:HA2	1:G:165:ILE:CD1	2.41	0.51
4:F:29:GLN:HG2	4:F:56:VAL:HB	1.93	0.51
4:F:115:LYS:NZ	7:F:404[B]:CL:CL	2.79	0.51
4:H:57:VAL:HG13	4:H:81:LEU:HD23	1.93	0.50
3:C:20:ILE:N	3:C:20:ILE:CD1	2.74	0.50
4:D:74:LYS:NZ	4:D:101:ASP:OD2	2.43	0.50
1:G:144:LEU:O	1:G:148:GLY:N	2.41	0.50
2:B:106:ASN:ND2	2:B:162:HIS:CE1	2.80	0.50
4:H:49:LEU:HD21	4:H:242:PRO:HB2	1.93	0.50
1:A:139:LEU:O	2:B:78:ASN:ND2	2.37	0.50
3:C:73:ILE:O	3:C:77:ILE:HG12	2.11	0.50
3:C:107:PRO:CD	3:C:132:ILE:HD12	2.41	0.50
3:C:181:SER:HA	3:C:206:THR:HB	1.94	0.50
4:H:244:GLU:HA	4:H:247:LYS:HD2	1.94	0.50
1:A:209:GLY:HA2	2:B:247:LYS:HD2	1.94	0.50
4:D:100:ILE:HG21	4:D:103:VAL:HG22	1.94	0.50
1:A:48[B]:ARG:NH2	8:A:556:HOH:O	2.40	0.50
1:A:58:ARG:NH2	6:A:401:SO4:O4	2.45	0.50
1:A:203[A]:LEU:HD12	1:A:233:THR:O	2.12	0.49
4:F:166:LEU:HD13	4:F:174:ILE:HD11	1.94	0.49
3:C:85:ILE:O	3:C:105:ILE:HD12	2.12	0.49
4:F:203:LEU:HD22	4:F:225:ILE:HG21	1.93	0.49
4:H:159:ALA:O	4:H:163:ILE:HD13	2.12	0.49
3:C:78:ASN:OD1	8:C:620:HOH:O	2.20	0.49
4:D:52:ALA:CB	4:D:243:VAL:HG13	2.42	0.49
4:H:256:PHE:N	4:H:257:GLY:HA2	2.28	0.49
1:G:112:ARG:NE	6:G:402:SO4:S	2.84	0.49
4:H:58:ARG:HD3	8:H:583:HOH:O	2.11	0.49
2:B:202:HIS:ND1	2:B:233:THR:HG23	2.27	0.49
4:F:87:PHE:CG	4:F:88:ASP:N	2.81	0.49
1:G:230:ILE:HG22	1:G:231:GLY:N	2.28	0.49
4:H:237:SER:C	4:H:238:LEU:HD12	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:E:536:HOH:O	4:H:217:LYS:HD2	2.11	0.49
1:A:181:SER:O	1:A:217:LYS:HE3	2.13	0.49
3:C:31:MSE:SE	3:C:58:ARG:NH2	2.96	0.48
5:E:144:LEU:C	8:E:647:HOH:O	2.56	0.48
2:B:216:VAL:HG11	3:C:251:GLU:HB3	1.95	0.48
1:G:96:ILE:HA	1:G:100:ILE:HG22	1.93	0.48
1:A:203[A]:LEU:HB3	1:A:222:LEU:HD22	1.94	0.48
4:F:101:ASP:O	4:F:128[A]:ILE:HG23	2.13	0.48
1:G:6:HIS:H	1:G:9:LYS:CE	2.27	0.48
1:A:110:ILE:HB	1:A:111:GLY:HA2	1.94	0.48
2:B:84:ASP:HA	2:B:104:ARG:HB3	1.95	0.48
1:G:6:HIS:HB3	1:G:9:LYS:HE2	1.95	0.48
1:A:88[A]:ASP:C	1:A:88[A]:ASP:OD1	2.56	0.48
4:D:188:GLU:O	4:D:192:LYS:HB2	2.14	0.48
4:D:139:LEU:HB3	4:D:143:ILE:HD11	1.95	0.48
1:G:21:GLY:HA2	1:G:232:ASN:ND2	2.29	0.47
5:E:203:LEU:HB3	5:E:222:LEU:HD22	1.96	0.47
1:A:169:LEU:HG	8:A:588:HOH:O	2.14	0.47
1:A:234:LEU:HD11	1:A:249:ALA:HB1	1.96	0.47
2:B:89:TYR:HD1	8:B:547:HOH:O	1.96	0.47
4:D:85:ILE:HD12	4:D:85:ILE:N	2.29	0.47
5:E:28:ILE:HD12	5:E:249:ALA:CB	2.44	0.47
1:A:8:THR:C	1:A:9:LYS:HD2	2.39	0.47
5:E:31:MSE:SE	5:E:58:ARG:NH2	2.98	0.47
1:A:203[B]:LEU:HB3	1:A:222:LEU:HD22	1.97	0.47
2:B:113:ARG:H	2:B:113:ARG:HD2	1.80	0.47
4:D:132:ILE:HD12	4:D:163:ILE:HD13	1.96	0.47
4:F:236:ILE:HG13	4:F:248:VAL:HG13	1.96	0.47
1:G:6:HIS:CB	1:G:9:LYS:HE2	2.45	0.47
1:A:97[B]:GLU:OE1	1:A:97[B]:GLU:HA	2.14	0.47
3:C:31:MSE:SE	3:C:58:ARG:CZ	3.13	0.47
3:C:101:ASP:O	3:C:129:PRO:HD2	2.15	0.47
4:D:159:ALA:O	4:D:163:ILE:HG12	2.14	0.47
5:E:46:ILE:HD12	5:E:57:VAL:HG21	1.96	0.47
5:E:142[B]:HIS:HE1	8:E:600:HOH:O	1.95	0.47
4:F:234:LEU:C	4:F:234:LEU:CD1	2.88	0.47
4:H:26:LEU:N	4:H:26:LEU:HD12	2.30	0.47
4:F:156:VAL:HG21	4:F:192:LYS:HB3	1.96	0.47
4:F:190:TYR:HD1	4:F:201:LEU:HD23	1.80	0.47
4:D:101:ASP:O	4:D:128:ILE:HG23	2.14	0.46
4:F:110:ILE:O	4:F:115:LYS:HE2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:71:ALA:O	4:H:75:LYS:HG3	2.15	0.46
1:A:142:HIS:HA	1:A:145[A]:GLU:HG2	1.97	0.46
1:G:228:LYS:O	1:G:228:LYS:HG3	2.15	0.46
1:A:143:ILE:N	1:A:143:ILE:HD12	2.31	0.46
1:A:164:LYS:HE2	1:A:168:ASP:OD1	2.15	0.46
5:E:26:LEU:CD1	5:E:258:LEU:HD13	2.45	0.46
3:C:139:LEU:O	4:D:78:ASN:OD1	2.34	0.46
1:A:234:LEU:C	1:A:234:LEU:CD1	2.88	0.46
3:C:28:ILE:HD11	3:C:253:LEU:HD12	1.96	0.46
4:D:74:LYS:HA	4:D:77:ILE:HG12	1.98	0.46
1:G:7:ARG:HG3	1:G:229:GLY:O	2.16	0.46
2:B:132:ILE:HG21	2:B:162:HIS:HB2	1.97	0.46
4:D:104:ARG:O	4:D:104:ARG:HD3	2.16	0.46
4:D:58:ARG:HE	4:D:84:ASP:HB3	1.81	0.46
1:A:29:GLN:HG2	1:A:56:VAL:HB	1.98	0.45
2:B:217:KCX:CX	8:B:581:HOH:O	2.64	0.45
1:G:149:TYR:HB2	1:G:150:PRO:CD	2.47	0.45
4:H:254:LYS:O	4:H:257:GLY:CA	2.64	0.45
4:D:58:ARG:NE	4:D:84:ASP:HB3	2.31	0.45
1:G:20:ILE:CD1	1:G:27:ILE:HG13	2.46	0.45
4:D:85:ILE:HB	4:D:105:ILE:CG2	2.47	0.45
4:D:202:HIS:HA	4:D:233:THR:CG2	2.46	0.45
5:E:174:ILE:O	5:E:200:PRO:HD2	2.17	0.45
4:D:156:VAL:HG21	4:D:192:LYS:HB3	1.97	0.45
4:F:132:ILE:HD11	4:F:166:LEU:HD12	1.97	0.45
1:G:112:ARG:NE	6:G:402:SO4:O2	2.50	0.45
4:D:131:ARG:HA	4:D:175:ILE:O	2.17	0.45
1:A:128:ILE:HG23	1:A:129:PRO:HD2	1.98	0.45
1:A:132:ILE:CD1	1:A:163:ILE:HD13	2.47	0.45
3:C:42:THR:O	3:C:46:ILE:HG12	2.17	0.45
1:A:203[A]:LEU:HD12	1:A:203[A]:LEU:H	1.82	0.45
5:E:102:KCX:OQ2	5:E:131:ARG:NE	2.47	0.45
4:F:113:ARG:HB3	6:F:402:SO4:O4	2.17	0.45
4:F:167:GLU:HG2	4:F:172:HIS:CE1	2.52	0.45
4:F:230:ILE:HG22	4:F:231:GLY:N	2.33	0.44
4:F:106:ASN:HA	4:F:107:PRO:HD3	1.88	0.44
3:C:28:ILE:HD12	3:C:249:ALA:CB	2.47	0.44
4:F:222:LEU:CD1	4:F:236:ILE:HD12	2.47	0.44
3:C:202:HIS:HA	3:C:233:THR:HG23	1.99	0.44
4:H:89:TYR:HB3	4:H:110:ILE:HD11	2.00	0.44
4:H:254:LYS:O	4:H:257:GLY:HA3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:190:TYR:HB2	4:D:225:ILE:HD12	2.00	0.44
4:F:111:GLY:HA3	4:F:115:LYS:HE2	2.00	0.44
2:B:224:ALA:O	2:B:228:LYS:HG2	2.17	0.44
4:D:179:LYS:HE3	4:D:206:THR:HG22	1.98	0.44
5:E:106:ASN:HB3	5:E:162:HIS:CE1	2.53	0.44
1:A:105:ILE:HD11	1:A:110:ILE:CD1	2.48	0.44
3:C:107:PRO:HD2	3:C:132:ILE:HD12	1.99	0.44
4:H:58:ARG:HH21	4:H:82:VAL:HG11	1.81	0.44
4:F:131[B]:ARG:HD3	4:F:177:SER:OG	2.17	0.43
3:C:142:HIS:O	3:C:145:GLU:N	2.51	0.43
1:A:203[B]:LEU:HD22	1:A:225:ILE:HG21	2.00	0.43
1:G:236:ILE:HG22	1:G:238:LEU:HG	2.00	0.43
1:A:251:GLU:OE2	4:D:213:ALA:HB1	2.18	0.43
3:C:7:ARG:HB2	3:C:232:ASN:ND2	2.34	0.43
2:B:109:ASN:O	2:B:110:ILE:C	2.61	0.43
4:F:122:ALA:HA	4:F:126:ARG:HH21	1.83	0.43
4:F:250:ARG:CG	4:F:250:ARG:NH1	2.81	0.43
1:G:26:LEU:HD22	1:G:226:LEU:HD21	2.01	0.43
2:B:175:ILE:HD12	2:B:200:PRO:HB2	2.00	0.43
4:D:134:VAL:HG21	4:D:159:ALA:HB2	2.01	0.43
3:C:55:GLN:O	3:C:79:ILE:HB	2.18	0.43
1:G:207:GLU:HB3	1:G:237:SER:OG	2.18	0.43
2:B:207:GLU:HA	8:B:533:HOH:O	2.18	0.43
5:E:126:ARG:HB3	5:E:126:ARG:CZ	2.48	0.43
3:C:26:LEU:HD12	3:C:258:LEU:HD13	2.01	0.42
4:H:77:ILE:HD11	4:H:81:LEU:HG	1.99	0.42
4:F:61:VAL:HG12	4:F:61:VAL:O	2.18	0.42
4:H:195:ARG:NE	8:H:637:HOH:O	2.51	0.42
1:A:89[A]:TYR:HD2	1:A:90[A]:ARG:HG3	1.85	0.42
4:D:77:ILE:HD13	4:D:81:LEU:HD21	2.02	0.42
1:A:228[A]:LYS:HA	1:A:228[A]:LYS:CE	2.45	0.42
5:E:31:MSE:SE	5:E:58:ARG:CZ	3.17	0.42
2:B:105:ILE:HD13	2:B:119:VAL:HG11	2.00	0.42
5:E:143:ILE:CD1	5:E:158:SER:HB2	2.49	0.42
1:G:217:LYS:NZ	8:G:544:HOH:O	2.35	0.42
4:H:32:THR:HB	4:H:45:GLU:OE1	2.19	0.42
4:D:29:GLN:HG2	4:D:56:VAL:HB	2.02	0.42
4:D:74:LYS:HE2	4:D:81:LEU:HD12	2.01	0.42
4:D:84:ASP:C	4:D:85:ILE:HD12	2.44	0.42
4:F:212:PHE:CE2	1:G:251:GLU:HG3	2.55	0.42
1:A:260:SER:O	1:A:261:ASN:CB	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:87:PHE:CD1	3:C:87:PHE:C	2.97	0.42
4:D:84:ASP:HB3	4:D:104:ARG:HD2	2.01	0.42
1:A:4:MSE:SE	1:A:191:GLU:HB2	2.69	0.42
4:F:66:ALA:HB2	8:F:575[B]:HOH:O	2.20	0.42
4:H:107:PRO:HD2	4:H:132:ILE:HD12	2.02	0.42
4:H:163:ILE:HG22	4:H:167:GLU:OE2	2.20	0.42
1:A:47:LYS:HD2	8:A:623:HOH:O	2.20	0.42
4:F:180:ALA:CB	4:F:185:LEU:HD22	2.50	0.42
1:G:201:LEU:HD12	1:G:230:ILE:CG2	2.50	0.42
1:A:57:VAL:HG13	1:A:81:LEU:HD12	2.02	0.41
2:B:207:GLU:CD	2:B:239:SER:H	2.28	0.41
1:A:55:GLN:O	1:A:80:PRO:HD2	2.20	0.41
1:A:102:KCX:HA	1:A:128:ILE:CG2	2.50	0.41
3:C:107:PRO:HD2	3:C:162:HIS:HB3	2.02	0.41
3:C:107:PRO:HD3	3:C:132:ILE:CD1	2.50	0.41
4:D:15:VAL:HG12	4:D:20:ILE:HG12	2.02	0.41
4:D:74:LYS:NZ	4:D:99:GLY:O	2.38	0.41
3:C:26:LEU:HD21	3:C:231:GLY:C	2.45	0.41
4:D:54:CYS:HB2	4:D:246:VAL:HG21	2.03	0.41
5:E:179:LYS:HD2	6:E:401:SO4:S	2.61	0.41
5:E:236:ILE:HG12	5:E:249:ALA:HB2	2.01	0.41
4:F:96:ILE:HA	4:F:100:ILE:HG22	2.03	0.41
4:H:28:ILE:N	4:H:28:ILE:HD12	2.35	0.41
4:H:149:TYR:HB2	4:H:150:PRO:CD	2.51	0.41
3:C:179:LYS:HE2	3:C:206:THR:HA	2.01	0.41
4:H:8:THR:HG22	4:H:24:ASN:HB2	2.02	0.41
5:E:142[B]:HIS:CE1	8:E:600:HOH:O	2.71	0.41
1:A:23:ASN:OD1	1:A:23:ASN:C	2.63	0.41
5:E:141:ARG:O	5:E:145:GLU:HG2	2.20	0.41
4:H:179:LYS:NZ	4:H:204:GLY:HA3	2.35	0.41
1:A:234:LEU:HD12	1:A:234:LEU:O	2.19	0.41
3:C:26:LEU:HD11	3:C:226:LEU:HD22	2.02	0.41
4:H:82:VAL:HG22	4:H:102:LYS:HB3	2.03	0.41
1:A:89[B]:TYR:CE1	1:A:90[B]:ARG:HG3	2.55	0.41
1:A:234:LEU:HD11	1:A:249:ALA:CB	2.50	0.41
2:B:103:VAL:CG1	2:B:105:ILE:HG23	2.51	0.41
3:C:46:ILE:CD1	3:C:57:VAL:HG21	2.51	0.41
3:C:102:KCX:HA	3:C:128:ILE:HG13	2.02	0.41
4:D:24:ASN:O	4:D:258:LEU:HD22	2.21	0.41
4:D:187[B]:ILE:O	4:D:191:GLU:HG3	2.21	0.41
5:E:2:ASN:N	8:E:603:HOH:O	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:179:LYS:NZ	6:H:401:SO4:O2	2.54	0.41
4:D:250:ARG:O	4:D:254:LYS:HB3	2.21	0.40
1:A:80:PRO:HA	1:A:101:ASP:OD2	2.21	0.40
1:A:203[A]:LEU:HD11	1:A:231:GLY:CA	2.48	0.40
4:F:236:ILE:CG1	4:F:248:VAL:HG13	2.51	0.40
4:D:83:ALA:HB2	4:D:100:ILE:HG12	2.02	0.40
4:D:136:ALA:HB2	4:D:155:MSE:CE	2.52	0.40
4:F:179:LYS:HE2	4:F:206:THR:HA	2.03	0.40
2:B:15:VAL:CG2	2:B:20:ILE:HD12	2.46	0.40
5:E:167:GLU:HG2	5:E:172:HIS:CE1	2.56	0.40
1:A:89[B]:TYR:HB2	1:A:119:VAL:HG21	2.04	0.40
2:B:166:LEU:HB3	2:B:171:PHE:HB3	2.04	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	268/276 (97%)	260 (97%)	7 (3%)	1 (0%)	30	17
1	G	255/276 (92%)	246 (96%)	9 (4%)	0	100	100
2	B	258/276 (94%)	245 (95%)	11 (4%)	2 (1%)	16	6
3	C	252/276 (91%)	242 (96%)	10 (4%)	0	100	100
4	D	251/276 (91%)	241 (96%)	10 (4%)	0	100	100
4	F	265/276 (96%)	260 (98%)	4 (2%)	1 (0%)	30	17
4	H	254/276 (92%)	247 (97%)	7 (3%)	0	100	100
5	E	258/276 (94%)	251 (97%)	7 (3%)	0	100	100
All	All	2061/2208 (93%)	1992 (97%)	65 (3%)	4 (0%)	43	31

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	110	ILE
1	A	87	PHE
2	B	112	ARG
4	F	242	PRO

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	219/215 (102%)	215 (98%)	4 (2%)	51	40
1	G	206/215 (96%)	202 (98%)	4 (2%)	50	38
2	B	209/215 (97%)	206 (99%)	3 (1%)	59	49
3	C	205/214 (96%)	202 (98%)	3 (2%)	57	47
4	D	206/216 (95%)	202 (98%)	4 (2%)	50	38
4	F	214/216 (99%)	211 (99%)	3 (1%)	59	49
4	H	206/216 (95%)	205 (100%)	1 (0%)	81	77
5	E	209/214 (98%)	205 (98%)	4 (2%)	50	38
All	All	1674/1721 (97%)	1648 (98%)	26 (2%)	59	44

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	31[A]	MSE
1	A	31[B]	MSE
1	A	234	LEU
2	B	31	MSE
2	B	113	ARG
2	B	234	LEU
3	C	10	THR
3	C	128	ILE
3	C	191	GLU
4	D	9[A]	LYS
4	D	9[B]	LYS

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Mol	Chain	Res	Type
4	D	124	LYS
4	D	211	LEU
5	E	94[A]	LYS
5	E	94[B]	LYS
5	E	132	ILE
5	E	183	VAL
4	F	48	ARG
4	F	234	LEU
4	F	250	ARG
1	G	144	LEU
1	G	174	ILE
1	G	211	LEU
1	G	234	LEU
4	H	78	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	161	HIS
2	B	78	ASN
2	B	227	ASN
3	C	37	HIS
4	D	17	ASN
4	D	142	HIS
4	D	232	ASN
5	E	2	ASN
5	E	161	HIS
5	E	232	ASN
4	F	37	HIS
4	F	68	ASN
4	F	86	HIS
4	F	135	ASN
1	G	17	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

7 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	$\# Z > 2$
2	KCX	B	217	2	10,11,12	0.70	0	6,12,14	0.74	0
5	KCX	E	102	5	10,11,12	0.88	0	6,12,14	0.42	0
1	KCX	G	102	1	10,11,12	0.56	0	6,12,14	1.92	2 (33%)
3	KCX	C	247	3	10,11,12	0.56	0	6,12,14	0.69	0
3	KCX	C	102	3	10,11,12	0.72	0	6,12,14	0.70	0
1	KCX	A	102	1	10,11,12	0.64	0	6,12,14	1.35	1 (16%)
5	KCX	E	35	5	10,11,12	0.81	0	6,12,14	1.12	1 (16%)

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
2	KCX	B	217	2	-	1/9/10/12	-
5	KCX	E	102	5	-	1/9/10/12	-
1	KCX	G	102	1	-	0/9/10/12	-
3	KCX	C	247	3	-	2/9/10/12	-
3	KCX	C	102	3	-	1/9/10/12	-
1	KCX	A	102	1	-	0/9/10/12	-
5	KCX	E	35	5	-	1/9/10/12	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	102	KCX	OQ1-CX-NZ	-3.75	119.23	124.92
1	A	102	KCX	OQ1-CX-NZ	-2.83	120.61	124.92
1	G	102	KCX	CE-NZ-CX	2.64	126.46	121.98
5	E	35	KCX	CE-NZ-CX	2.13	125.60	121.98

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	102	KCX	O-C-CA-CB
5	E	102	KCX	O-C-CA-CB
2	B	217	KCX	CG-CD-CE-NZ
5	E	35	KCX	CG-CD-CE-NZ
3	C	247	KCX	CE-CD-CG-CB
3	C	247	KCX	C-CA-CB-CG

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	217	KCX	1	0
5	E	102	KCX	1	0
1	G	102	KCX	1	0
3	C	102	KCX	2	0
1	A	102	KCX	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 23 ligands modelled in this entry, 10 are monoatomic - leaving 13 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	$\# Z > 2$
6	SO4	B	401	-	4,4,4	0.43	0	6,6,6	0.25	0
6	SO4	C	401	-	4,4,4	0.36	0	6,6,6	0.25	0
6	SO4	E	402	-	4,4,4	0.41	0	6,6,6	0.13	0
6	SO4	A	401	-	4,4,4	0.42	0	6,6,6	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	D	402	-	4,4,4	0.40	0	6,6,6	0.08	0
6	SO4	E	401	-	4,4,4	0.40	0	6,6,6	0.41	0
6	SO4	H	401	-	4,4,4	0.49	0	6,6,6	0.18	0
6	SO4	D	401	-	4,4,4	0.44	0	6,6,6	0.21	0
6	SO4	F	401	-	4,4,4	0.39	0	6,6,6	0.30	0
6	SO4	G	402	-	4,4,4	0.43	0	6,6,6	0.08	0
6	SO4	F	402	-	4,4,4	0.45	0	6,6,6	0.23	0
6	SO4	G	401	-	4,4,4	0.34	0	6,6,6	0.41	0
6	SO4	C	402	-	4,4,4	0.50	0	6,6,6	0.17	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.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	401	SO4	1	0
6	A	401	SO4	1	0
6	E	401	SO4	2	0
6	H	401	SO4	2	0
6	G	402	SO4	2	0
6	F	402	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	255/276 (92%)	-1.14	0 100 100	7, 21, 33, 42	9 (3%)
1	G	253/276 (91%)	-1.23	0 100 100	11, 17, 26, 32	0
2	B	254/276 (92%)	-1.10	0 100 100	9, 24, 40, 49	2 (0%)
3	C	251/276 (90%)	-1.11	0 100 100	13, 22, 32, 54	1 (0%)
4	D	247/276 (89%)	-1.06	0 100 100	9, 24, 41, 51	4 (1%)
4	F	258/276 (93%)	-1.16	0 100 100	9, 19, 29, 43	5 (1%)
4	H	251/276 (90%)	-1.19	0 100 100	9, 19, 29, 37	1 (0%)
5	E	254/276 (92%)	-1.24	0 100 100	9, 17, 24, 30	2 (0%)
All	All	2023/2208 (91%)	-1.15	0 100 100	7, 20, 33, 54	24 (1%)

There are no RSRZ outliers to report.

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	KCX	B	217	12/13	0.98	0.04	18,24,36,37	0
3	KCX	C	247	12/13	0.98	0.04	16,23,32,35	0
5	KCX	E	35	12/13	0.98	0.04	15,17,27,28	0
1	KCX	G	102	12/13	0.98	0.05	16,20,28,32	0
3	KCX	C	102	12/13	0.99	0.04	17,18,19,20	0
5	KCX	E	102	12/13	0.99	0.03	10,11,13,14	0
1	KCX	A	102	12/13	0.99	0.04	15,16,18,18	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
6	SO4	G	402	5/5	0.98	0.07	34,43,45,46	0
6	SO4	C	401	5/5	0.99	0.04	25,26,27,27	0
6	SO4	C	402	5/5	0.99	0.04	29,31,35,38	0
6	SO4	D	401	5/5	0.99	0.04	32,34,37,38	0
6	SO4	D	402	5/5	0.99	0.02	39,39,39,40	5
6	SO4	E	402	5/5	0.99	0.06	38,42,43,45	0
6	SO4	F	401	5/5	0.99	0.04	33,36,41,42	0
6	SO4	F	402	5/5	0.99	0.03	37,37,39,39	0
6	SO4	G	401	5/5	0.99	0.03	16,19,20,21	0
6	SO4	B	401	5/5	0.99	0.05	29,31,35,35	0
6	SO4	H	401	5/5	0.99	0.04	25,30,31,36	0
7	CL	A	402	1/1	0.99	0.08	53,53,53,53	0
7	CL	B	402[A]	1/1	0.99	0.06	32,32,32,32	1
7	CL	B	402[B]	1/1	0.99	0.06	21,21,21,21	1
7	CL	C	403	1/1	0.99	0.05	37,37,37,37	0
7	CL	C	404	1/1	0.99	0.04	44,44,44,44	0
7	CL	C	405	1/1	0.99	0.07	50,50,50,50	0
7	CL	F	404[A]	1/1	0.99	0.08	31,31,31,31	1
7	CL	F	404[B]	1/1	0.99	0.08	23,23,23,23	1
7	CL	B	403	1/1	1.00	0.07	42,42,42,42	1
7	CL	F	403	1/1	1.00	0.04	30,30,30,30	0
6	SO4	E	401	5/5	1.00	0.03	18,19,21,21	0
6	SO4	A	401	5/5	1.00	0.04	18,21,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.