



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:45 PM UTC

PDB ID : 3GTS / pdb_00003gts
Title : Crystal Structure of Dicamba Monooxygenase with Non-heme Iron and Dicamba
Authors : Rydel, T.J.; Sturman, E.J.; Moshiri, F.; Brown, G.R.; Qi, Y.
Deposited on : 2009-03-28
Resolution : 2.20 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

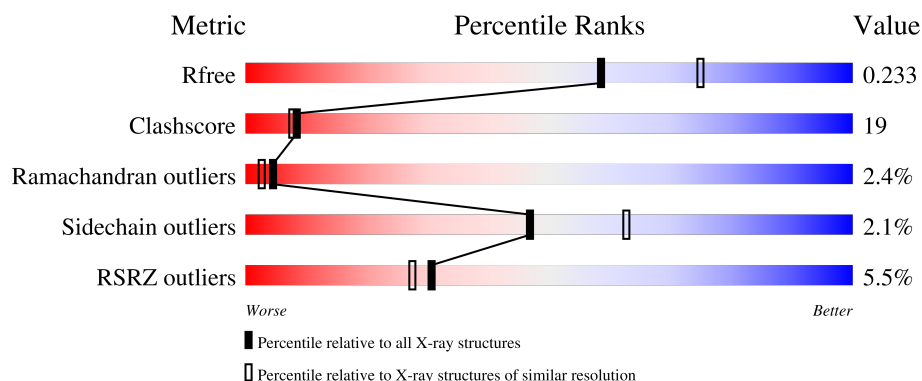
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	349	0% 67% 28% 5%
1	B	349	6% 63% 32% 5%
1	C	349	8% 61% 27% 5% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8329 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 DdmC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	B	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	C	327	Total	C	N	O	S	0	0	0
			2524	1594	450	467	13			

There are 33 discrepancies between the modelled and reference sequences:

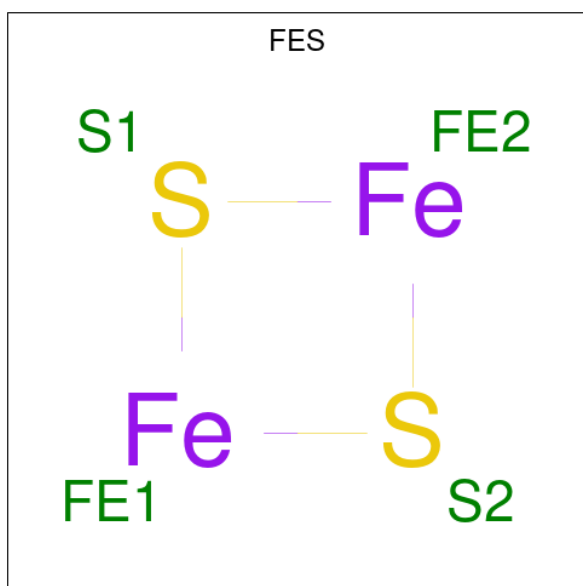
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q5S3I3
A	2	ALA	MET	engineered mutation	UNP Q5S3I3
A	341	ARG	-	expression tag	UNP Q5S3I3
A	342	LEU	-	expression tag	UNP Q5S3I3
A	343	GLU	-	expression tag	UNP Q5S3I3
A	344	HIS	-	expression tag	UNP Q5S3I3
A	345	HIS	-	expression tag	UNP Q5S3I3
A	346	HIS	-	expression tag	UNP Q5S3I3
A	347	HIS	-	expression tag	UNP Q5S3I3
A	348	HIS	-	expression tag	UNP Q5S3I3
A	349	HIS	-	expression tag	UNP Q5S3I3
B	2	ALA	-	expression tag	UNP Q5S3I3
B	2	ALA	MET	engineered mutation	UNP Q5S3I3
B	341	ARG	-	expression tag	UNP Q5S3I3
B	342	LEU	-	expression tag	UNP Q5S3I3
B	343	GLU	-	expression tag	UNP Q5S3I3
B	344	HIS	-	expression tag	UNP Q5S3I3
B	345	HIS	-	expression tag	UNP Q5S3I3
B	346	HIS	-	expression tag	UNP Q5S3I3
B	347	HIS	-	expression tag	UNP Q5S3I3
B	348	HIS	-	expression tag	UNP Q5S3I3
B	349	HIS	-	expression tag	UNP Q5S3I3
C	2	ALA	-	expression tag	UNP Q5S3I3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2	ALA	MET	engineered mutation	UNP Q5S3I3
C	341	ARG	-	expression tag	UNP Q5S3I3
C	342	LEU	-	expression tag	UNP Q5S3I3
C	343	GLU	-	expression tag	UNP Q5S3I3
C	344	HIS	-	expression tag	UNP Q5S3I3
C	345	HIS	-	expression tag	UNP Q5S3I3
C	346	HIS	-	expression tag	UNP Q5S3I3
C	347	HIS	-	expression tag	UNP Q5S3I3
C	348	HIS	-	expression tag	UNP Q5S3I3
C	349	HIS	-	expression tag	UNP Q5S3I3

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe S 4 2 2	0	0
2	B	1	Total Fe S 4 2 2	0	0
2	C	1	Total Fe S 4 2 2	0	0

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

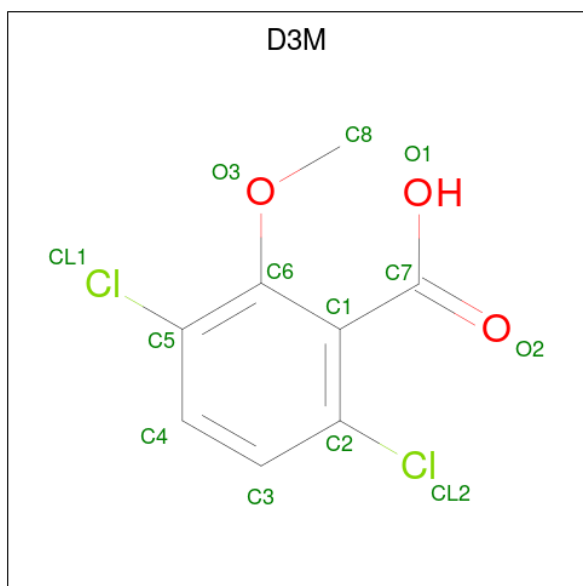
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Fe	0	0
			1	1		

- Molecule 4 is 3,6-dichloro-2-methoxybenzoic acid (CCD ID: D3M) (formula: $C_8H_6Cl_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Cl	O	0	0
			13	8	2	3		
4	B	1	Total	C	Cl	O	0	0
			13	8	2	3		

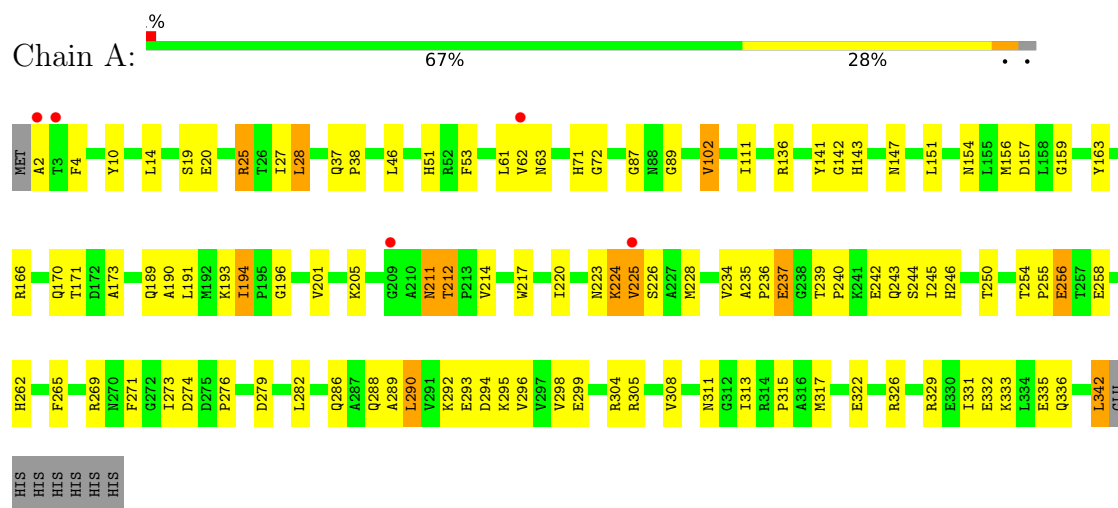
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	169	Total	O	0	0
			169	169		
5	B	155	Total	O	0	0
			155	155		
5	C	159	Total	O	0	0
			159	159		

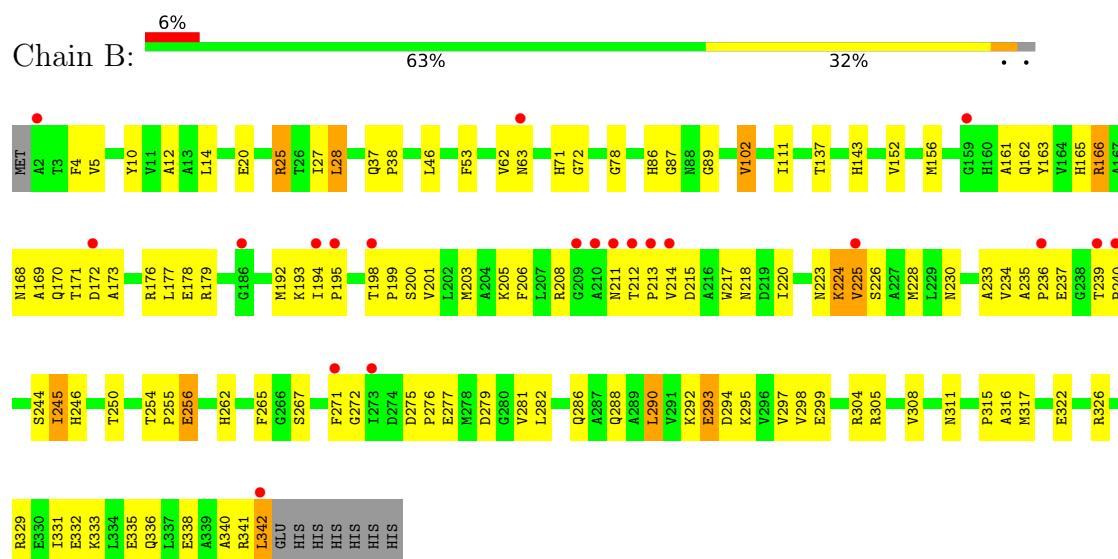
3 Residue-property plots

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 and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: DdmC

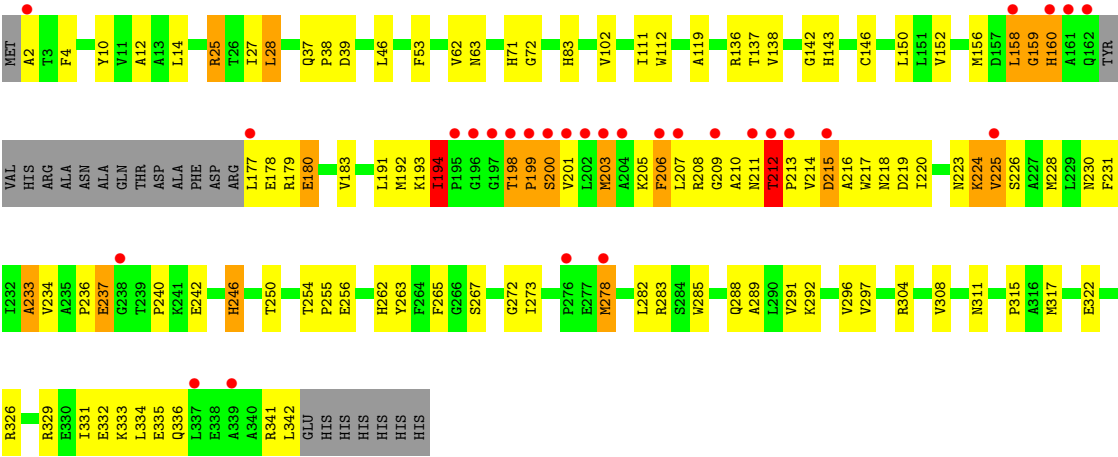


• Molecule 1: DdmC



• Molecule 1: DdmC





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	80.80Å 80.80Å 159.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.20 20.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	84.9 (20.00-2.20) 85.0 (20.00-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 2.20Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.230 , 0.266 0.205 , 0.233	Depositor DCC
R_{free} test set	5061 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.208	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.003 for -h,-k,l 0.046 for h,-h-k,-l 0.022 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8329	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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: FES, FE, D3M

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	0.39	0/2706	0.94	11/3685 (0.3%)
1	B	0.38	0/2706	0.94	14/3685 (0.4%)
1	C	0.40	0/2585	0.94	14/3519 (0.4%)
All	All	0.39	0/7997	0.94	39/10889 (0.4%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	194	ILE	N-CA-C	7.72	115.79	108.15
1	A	72	GLY	N-CA-C	7.49	125.71	115.30
1	B	72	GLY	N-CA-C	7.49	125.71	115.30
1	C	72	GLY	N-CA-C	7.38	125.56	115.30
1	A	226	SER	N-CA-C	7.06	118.74	110.41
1	C	246	HIS	N-CA-C	6.91	118.60	108.86
1	C	226	SER	N-CA-C	6.88	118.53	110.41
1	B	226	SER	N-CA-C	6.74	118.36	110.41
1	B	290	LEU	N-CA-C	6.65	118.32	111.14
1	A	194	ILE	CA-C-N	6.15	126.06	119.85
1	A	194	ILE	C-N-CA	6.15	126.06	119.85
1	B	14	LEU	N-CA-C	-6.04	101.95	110.29
1	C	278	MET	N-CA-C	6.04	118.66	111.71
1	A	246	HIS	N-CA-C	5.87	117.17	108.60
1	B	246	HIS	N-CA-C	5.87	117.71	108.96
1	C	46	LEU	N-CA-C	-5.85	101.19	109.96
1	B	293	GLU	N-CA-C	5.82	117.29	111.07
1	C	297	VAL	N-CA-C	5.77	115.95	110.53
1	A	14	LEU	N-CA-C	-5.74	102.37	110.29
1	A	46	LEU	N-CA-C	-5.70	101.61	110.10
1	C	152	VAL	N-CA-C	-5.64	105.23	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	311	ASN	N-CA-C	-5.63	101.64	110.36
1	B	46	LEU	N-CA-C	-5.61	101.74	110.10
1	C	14	LEU	N-CA-C	-5.48	102.73	110.29
1	C	233	ALA	N-CA-C	5.39	116.99	108.96
1	B	176	ARG	N-CA-C	-5.38	106.03	112.59
1	A	256	GLU	N-CA-C	-5.34	104.43	111.96
1	C	53	PHE	N-CA-C	5.32	119.72	113.23
1	A	53	PHE	N-CA-C	5.25	119.64	113.23
1	A	313	ILE	N-CA-C	5.24	115.85	109.30
1	A	311	ASN	N-CA-C	-5.22	102.27	110.36
1	B	53	PHE	N-CA-C	5.14	119.50	113.23
1	C	311	ASN	N-CA-C	-5.14	102.40	110.36
1	B	237	GLU	CB-CA-C	-5.12	110.26	117.23
1	C	194	ILE	CA-C-N	5.11	125.03	119.76
1	C	194	ILE	C-N-CA	5.11	125.03	119.76
1	B	256	GLU	N-CA-C	-5.08	104.80	111.96
1	B	297	VAL	N-CA-C	5.06	115.28	110.42
1	B	152	VAL	N-CA-C	-5.04	105.79	110.53

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	2641	0	2577	85	0
1	B	2641	0	2577	109	0
1	C	2524	0	2472	104	0
2	A	4	0	0	1	0
2	B	4	0	0	1	0
2	C	4	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	13	0	5	0	0
4	B	13	0	5	0	0
5	A	169	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	155	0	0	7	0
5	C	159	0	0	9	0
All	All	8329	0	7636	295	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 19.

All (295) 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:A:28:LEU:HD22	1:A:255:PRO:HG2	1.41	1.02
1:C:156:MET:HE3	1:C:228:MET:SD	2.05	0.97
1:B:28:LEU:HD22	1:B:255:PRO:HG2	1.48	0.94
1:C:28:LEU:HD22	1:C:255:PRO:HG2	1.50	0.93
1:A:245:ILE:HD12	1:A:271:PHE:HB3	1.50	0.92
1:B:156:MET:HE3	1:B:228:MET:SD	2.11	0.90
1:B:171:THR:HG22	1:B:173:ALA:H	1.37	0.88
1:A:156:MET:HE3	1:A:228:MET:SD	2.15	0.86
1:C:272:GLY:HA3	1:C:278:MET:HE1	1.58	0.85
1:B:286:GLN:HA	1:B:290:LEU:HD23	1.57	0.85
1:A:171:THR:HG22	1:A:173:ALA:H	1.39	0.84
1:B:203:MET:HE3	1:B:234:VAL:HB	1.57	0.83
1:C:342:LEU:HD23	1:C:342:LEU:H	1.44	0.82
1:B:171:THR:HG21	1:B:194:ILE:HG21	1.62	0.79
1:B:211:ASN:ND2	1:B:212:THR:H	1.81	0.78
1:A:286:GLN:HA	1:A:290:LEU:HD23	1.67	0.77
1:C:233:ALA:HB2	1:C:246:HIS:HB3	1.65	0.77
1:A:295:LYS:O	1:A:299:GLU:HG3	1.84	0.77
1:C:194:ILE:HG13	1:C:216:ALA:HB3	1.67	0.76
1:B:217:TRP:HB2	1:B:233:ALA:HB3	1.68	0.75
1:A:245:ILE:HG23	1:A:271:PHE:CD2	2.21	0.75
1:B:156:MET:HE1	1:B:220:ILE:HG12	1.69	0.74
1:A:156:MET:HE1	1:A:220:ILE:HG12	1.67	0.74
1:C:177:LEU:HD12	1:C:178:GLU:H	1.53	0.73
1:B:63:ASN:HB3	5:B:1001:HOH:O	1.88	0.73
1:B:192:MET:HE2	1:B:218:ASN:HB2	1.70	0.73
1:A:290:LEU:HD22	1:A:290:LEU:H	1.53	0.72
1:A:143:HIS:HE1	1:A:256:GLU:OE1	1.72	0.72
1:A:342:LEU:HD23	1:A:342:LEU:H	1.54	0.72
1:A:191:LEU:N	1:A:191:LEU:HD12	2.05	0.71
1:A:214:VAL:HG21	1:A:234:VAL:HG22	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:ARG:HB3	1:B:342:LEU:HD23	1.73	0.70
1:B:342:LEU:HD23	1:B:342:LEU:H	1.57	0.70
1:B:214:VAL:HG21	1:B:234:VAL:HG22	1.74	0.69
1:C:288:GLN:HE21	1:C:288:GLN:HA	1.59	0.68
1:B:245:ILE:HG12	1:B:271:PHE:HB3	1.75	0.68
1:B:295:LYS:O	1:B:299:GLU:HG3	1.93	0.68
1:A:156:MET:HE2	1:A:156:MET:HA	1.76	0.67
1:C:203:MET:HE2	1:C:203:MET:HA	1.76	0.67
1:A:290:LEU:HD22	1:A:290:LEU:N	2.10	0.66
1:B:178:GLU:OE2	1:B:193:LYS:HE2	1.96	0.65
1:C:4:PHE:HB3	1:C:28:LEU:HD13	1.79	0.65
1:B:177:LEU:HD11	1:B:192:MET:SD	2.38	0.64
1:C:143:HIS:HE1	1:C:256:GLU:OE1	1.80	0.63
1:B:143:HIS:HE1	1:B:256:GLU:OE1	1.82	0.62
1:C:198:THR:H	1:C:199:PRO:HD2	1.64	0.62
1:C:193:LYS:HE3	1:C:217:TRP:HE1	1.65	0.62
1:B:331:ILE:O	1:B:335:GLU:HG3	2.00	0.62
1:B:245:ILE:HG23	1:B:271:PHE:CD2	2.35	0.61
1:B:192:MET:HE3	1:B:194:ILE:HD11	1.82	0.61
1:C:212:THR:O	1:C:214:VAL:HG13	2.01	0.61
1:B:267:SER:OG	1:B:282:LEU:HD13	2.01	0.61
1:C:288:GLN:HB2	5:C:988:HOH:O	2.00	0.61
1:B:4:PHE:HB3	1:B:28:LEU:HD13	1.83	0.60
1:B:156:MET:HE2	1:B:156:MET:HA	1.83	0.60
1:B:172:ASP:OD1	1:B:198:THR:HB	2.00	0.60
1:C:37:GLN:HB3	1:C:38:PRO:HD2	1.83	0.60
1:B:208:ARG:HH22	1:B:272:GLY:HA2	1.67	0.60
1:C:156:MET:HE2	1:C:156:MET:HA	1.83	0.60
1:C:233:ALA:CB	1:C:246:HIS:HB3	2.32	0.60
1:C:315:PRO:HB2	1:C:317:MET:HE3	1.84	0.60
1:B:71:HIS:HB2	2:B:501:FES:S1	2.43	0.59
1:C:331:ILE:O	1:C:335:GLU:HG3	2.02	0.59
1:C:207:LEU:C	1:C:209:GLY:H	2.11	0.59
1:A:19:SER:HB2	5:A:853:HOH:O	2.03	0.59
1:A:315:PRO:HB2	1:A:317:MET:HE3	1.84	0.59
1:B:27:ILE:HD12	1:B:254:THR:HG21	1.85	0.59
1:B:315:PRO:HB2	1:B:317:MET:HE3	1.84	0.58
1:C:156:MET:HE1	1:C:220:ILE:HG12	1.83	0.58
1:A:37:GLN:HB3	1:A:38:PRO:HD2	1.84	0.58
1:B:162:GLN:HE22	1:B:171:THR:H	1.50	0.58
1:C:211:ASN:O	1:C:212:THR:HG22	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ILE:O	1:A:335:GLU:HG3	2.03	0.58
1:B:37:GLN:HB3	1:B:38:PRO:HD2	1.85	0.57
1:C:193:LYS:HG2	1:C:217:TRP:CD1	2.39	0.57
1:B:28:LEU:CD2	1:B:255:PRO:HG2	2.30	0.57
1:B:165:HIS:HA	1:B:293:GLU:OE2	2.04	0.57
1:C:265:PHE:CE2	1:C:285:TRP:HE3	2.23	0.56
1:C:137:THR:HA	1:C:267:SER:O	2.05	0.56
1:B:10:TYR:CE2	1:B:225:VAL:HG11	2.41	0.56
1:B:20:GLU:HB2	5:B:815:HOH:O	2.04	0.56
1:A:4:PHE:HB3	1:A:28:LEU:HD13	1.86	0.56
1:B:214:VAL:HG23	1:B:235:ALA:O	2.05	0.56
1:C:206:PHE:HE2	1:C:246:HIS:HA	1.71	0.56
1:B:208:ARG:HH22	1:B:272:GLY:CA	2.20	0.55
1:C:206:PHE:HD1	1:C:206:PHE:H	1.54	0.55
1:A:242:GLU:CD	1:A:242:GLU:H	2.15	0.55
1:B:276:PRO:O	1:B:279:ASP:HB2	2.06	0.55
1:A:147:ASN:ND2	5:A:890:HOH:O	2.40	0.55
1:C:177:LEU:O	1:C:193:LYS:HB2	2.06	0.55
1:C:199:PRO:HG2	1:C:200:SER:H	1.72	0.55
1:C:71:HIS:HB2	2:C:501:FES:S1	2.47	0.54
1:A:71:HIS:HB2	2:A:501:FES:S1	2.47	0.54
1:C:236:PRO:O	1:C:237:GLU:C	2.51	0.54
1:A:211:ASN:O	1:A:212:THR:HB	2.07	0.54
1:B:201:VAL:HG22	1:B:205:LYS:HE3	1.88	0.54
1:B:235:ALA:HB1	1:B:239:THR:OG1	2.08	0.54
1:A:102:VAL:HB	1:A:111:ILE:HD13	1.89	0.54
1:B:177:LEU:HD11	1:B:192:MET:CG	2.37	0.54
1:B:236:PRO:O	1:B:239:THR:HG23	2.08	0.54
1:C:233:ALA:HB2	1:C:246:HIS:CB	2.36	0.54
1:B:342:LEU:HD23	1:B:342:LEU:N	2.22	0.54
1:B:25:ARG:HD2	1:B:262:HIS:CE1	2.43	0.53
1:C:283:ARG:HG2	5:C:1164:HOH:O	2.06	0.53
1:A:240:PRO:HG2	1:A:243:GLN:HB2	1.90	0.53
1:A:239:THR:HB	1:A:240:PRO:HD2	1.91	0.53
1:B:193:LYS:HG2	1:B:195:PRO:HD3	1.90	0.53
1:C:317:MET:HA	1:C:317:MET:HE2	1.91	0.53
1:C:207:LEU:C	1:C:207:LEU:HD23	2.34	0.53
1:C:207:LEU:O	1:C:209:GLY:N	2.41	0.53
1:A:25:ARG:HD2	1:A:262:HIS:CE1	2.44	0.52
1:A:282:LEU:O	1:A:286:GLN:HG3	2.09	0.52
1:B:322:GLU:HG2	1:B:326:ARG:NH1	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:VAL:HG12	1:A:63:ASN:ND2	2.25	0.52
1:A:240:PRO:HG3	1:A:243:GLN:OE1	2.10	0.52
1:B:203:MET:O	1:B:206:PHE:HB2	2.09	0.52
1:A:224:LYS:HB2	5:A:894:HOH:O	2.09	0.52
1:A:269:ARG:HD3	1:A:282:LEU:HD11	1.91	0.52
1:A:290:LEU:H	1:A:290:LEU:CD2	2.20	0.52
1:A:294:ASP:O	1:A:298:VAL:HG22	2.10	0.52
1:C:214:VAL:HG12	1:C:236:PRO:HA	1.92	0.52
1:C:178:GLU:HG3	1:C:179:ARG:H	1.76	0.52
1:B:277:GLU:O	1:B:281:VAL:HG23	2.10	0.51
1:B:28:LEU:N	1:B:28:LEU:HD23	2.25	0.51
1:C:213:PRO:O	1:C:236:PRO:HA	2.10	0.51
1:A:166:ARG:HB2	1:A:170:GLN:HE21	1.76	0.51
1:B:317:MET:HE2	1:B:317:MET:HA	1.93	0.51
1:C:179:ARG:O	1:C:180:GLU:HB2	2.09	0.51
1:B:163:TYR:O	1:B:166:ARG:HB2	2.12	0.50
1:B:288:GLN:HE22	1:B:292:LYS:HD2	1.75	0.50
1:A:326:ARG:HG3	1:A:326:ARG:HH11	1.76	0.50
1:C:25:ARG:HD3	5:C:1152:HOH:O	2.09	0.50
1:A:245:ILE:HD12	1:A:271:PHE:CB	2.32	0.50
1:B:71:HIS:HD2	1:C:160:HIS:CD2	2.30	0.50
1:C:206:PHE:CD1	1:C:206:PHE:N	2.79	0.50
1:A:289:ALA:HB3	1:A:290:LEU:HD22	1.94	0.50
1:A:10:TYR:CZ	1:A:225:VAL:HG11	2.47	0.50
1:A:191:LEU:N	1:A:191:LEU:CD1	2.74	0.50
1:B:179:ARG:HB3	5:B:1027:HOH:O	2.12	0.50
1:C:192:MET:HE2	1:C:218:ASN:HD22	1.76	0.50
1:A:317:MET:HE2	1:A:317:MET:HA	1.94	0.50
1:B:201:VAL:O	1:B:205:LYS:HG3	2.11	0.50
1:A:322:GLU:HG2	1:A:326:ARG:NH1	2.27	0.49
1:C:288:GLN:NE2	5:C:988:HOH:O	2.46	0.49
1:A:304:ARG:HG3	1:A:304:ARG:HH11	1.77	0.49
1:B:137:THR:HA	1:B:267:SER:O	2.12	0.49
1:C:62:VAL:HG12	1:C:63:ASN:ND2	2.26	0.49
1:B:10:TYR:CZ	1:B:225:VAL:HG11	2.47	0.49
1:B:102:VAL:HB	1:B:111:ILE:HD13	1.93	0.49
1:B:173:ALA:HB3	1:B:194:ILE:CG2	2.43	0.49
1:C:28:LEU:HD23	1:C:28:LEU:N	2.27	0.49
1:B:20:GLU:HG3	5:B:997:HOH:O	2.12	0.49
1:A:274:ASP:HB3	5:A:710:HOH:O	2.12	0.49
1:B:214:VAL:HG22	1:B:215:ASP:N	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:VAL:HB	1:A:111:ILE:CD1	2.43	0.48
1:A:2:ALA:N	1:A:258:GLU:OE1	2.47	0.48
1:A:201:VAL:HG12	1:A:205:LYS:HE2	1.95	0.48
1:A:288:GLN:HE21	1:A:293:GLU:HG3	1.78	0.48
1:A:214:VAL:CG2	1:A:234:VAL:HG22	2.43	0.48
1:A:292:LYS:O	1:A:296:VAL:HG23	2.14	0.48
1:B:201:VAL:CG2	1:B:205:LYS:HE3	2.43	0.48
1:C:292:LYS:O	1:C:296:VAL:HG23	2.14	0.48
1:C:332:GLU:O	1:C:336:GLN:HG3	2.14	0.48
1:A:276:PRO:O	1:A:279:ASP:HB2	2.13	0.48
1:B:282:LEU:O	1:B:286:GLN:HG3	2.14	0.48
1:A:27:ILE:HG22	1:A:28:LEU:HD23	1.96	0.48
1:A:342:LEU:H	1:A:342:LEU:CD2	2.25	0.48
1:B:214:VAL:CG2	1:B:234:VAL:HG22	2.43	0.48
1:C:201:VAL:C	1:C:203:MET:H	2.20	0.48
1:C:206:PHE:CE2	1:C:246:HIS:HA	2.49	0.48
1:C:288:GLN:HE21	1:C:288:GLN:CA	2.22	0.48
1:A:193:LYS:HD2	1:A:217:TRP:HE1	1.79	0.47
1:B:62:VAL:HG12	1:B:63:ASN:ND2	2.28	0.47
1:B:304:ARG:HG3	1:B:304:ARG:HH11	1.79	0.47
1:A:163:TYR:HB3	1:A:166:ARG:HH21	1.79	0.47
1:C:178:GLU:HG3	1:C:179:ARG:N	2.29	0.47
1:A:51:HIS:HD2	1:B:316:ALA:O	1.98	0.47
1:B:5:VAL:HG13	1:B:338:GLU:OE2	2.14	0.47
1:B:340:ALA:O	1:B:341:ARG:C	2.57	0.47
1:C:198:THR:HB	1:C:199:PRO:HD3	1.96	0.47
1:A:235:ALA:HB2	1:A:244:SER:OG	2.14	0.47
1:B:78:GLY:HA3	5:B:872:HOH:O	2.14	0.47
1:B:326:ARG:HH11	1:B:326:ARG:HG3	1.80	0.47
1:C:10:TYR:CZ	1:C:225:VAL:HG11	2.50	0.47
1:A:156:MET:HE1	1:A:220:ILE:CG1	2.41	0.47
1:B:173:ALA:HB3	1:B:194:ILE:HG22	1.97	0.47
1:C:102:VAL:HB	1:C:111:ILE:HD13	1.96	0.47
1:C:205:LYS:HD3	1:C:205:LYS:H	1.80	0.47
1:C:288:GLN:HG2	5:C:802:HOH:O	2.15	0.47
1:C:322:GLU:HG2	1:C:326:ARG:NH1	2.29	0.47
1:B:168:ASN:O	1:B:200:SER:HB2	2.14	0.47
1:C:224:LYS:HG3	1:C:334:LEU:HD11	1.96	0.47
1:C:272:GLY:HA3	1:C:278:MET:CE	2.38	0.46
1:A:304:ARG:O	1:A:308:VAL:HG13	2.15	0.46
1:A:211:ASN:N	1:A:211:ASN:HD22	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:304:ARG:O	1:C:308:VAL:HG13	2.15	0.46
1:C:12:ALA:HB2	1:C:111:ILE:HG13	1.98	0.46
1:B:275:ASP:OD1	1:B:277:GLU:HB2	2.16	0.46
1:C:10:TYR:CE2	1:C:225:VAL:HG11	2.50	0.46
1:A:201:VAL:O	1:A:205:LYS:HG3	2.16	0.46
1:B:168:ASN:HA	1:B:201:VAL:CG1	2.46	0.46
1:B:290:LEU:H	1:B:290:LEU:HD22	1.80	0.46
1:B:288:GLN:NE2	1:B:292:LYS:HD2	2.31	0.46
1:C:158:LEU:O	1:C:160:HIS:N	2.49	0.46
1:A:28:LEU:HD23	1:A:28:LEU:N	2.31	0.45
1:C:158:LEU:O	1:C:159:GLY:C	2.59	0.45
1:B:102:VAL:HB	1:B:111:ILE:CD1	2.46	0.45
1:C:183:VAL:CG2	1:C:326:ARG:HD3	2.46	0.45
1:A:154:ASN:HA	5:C:703:HOH:O	2.17	0.45
1:B:290:LEU:HD22	1:B:290:LEU:N	2.32	0.45
1:B:211:ASN:ND2	1:B:212:THR:N	2.57	0.45
1:B:304:ARG:O	1:B:308:VAL:HG13	2.17	0.45
1:C:142:GLY:HA3	1:C:285:TRP:CZ3	2.51	0.45
1:C:206:PHE:HD2	1:C:234:VAL:HG12	1.82	0.45
1:A:288:GLN:HE21	1:A:293:GLU:CD	2.25	0.45
1:C:218:ASN:HA	1:C:231:PHE:O	2.18	0.45
1:C:83:HIS:HD2	5:C:952:HOH:O	1.99	0.44
1:A:342:LEU:HG	1:A:342:LEU:O	2.16	0.44
1:B:250:THR:O	1:B:265:PHE:HA	2.17	0.44
1:C:191:LEU:CD2	1:C:219:ASP:HB3	2.48	0.44
1:A:189:GLN:HG3	5:A:892:HOH:O	2.16	0.44
1:A:28:LEU:CD2	1:A:255:PRO:HG2	2.29	0.44
1:C:138:VAL:HG21	1:C:282:LEU:HB2	1.98	0.44
1:C:326:ARG:HG3	1:C:326:ARG:HH11	1.82	0.44
1:C:250:THR:O	1:C:265:PHE:HA	2.18	0.44
1:C:304:ARG:HH11	1:C:304:ARG:HG3	1.83	0.44
1:B:206:PHE:C	1:B:208:ARG:H	2.24	0.44
1:A:10:TYR:CE2	1:A:225:VAL:HG11	2.53	0.44
1:B:332:GLU:O	1:B:336:GLN:HG3	2.18	0.44
1:C:28:LEU:CD2	1:C:255:PRO:HG2	2.34	0.44
1:C:206:PHE:CD2	1:C:234:VAL:HG12	2.52	0.44
1:C:223:ASN:O	1:C:224:LYS:C	2.61	0.44
1:B:244:SER:HA	5:B:1056:HOH:O	2.17	0.43
1:B:288:GLN:O	1:B:292:LYS:HB3	2.18	0.43
1:C:205:LYS:HB2	1:C:205:LYS:HE2	1.80	0.43
1:C:25:ARG:HD2	1:C:262:HIS:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:CYS:HB2	1:C:150:LEU:HD12	2.01	0.43
1:A:27:ILE:HD12	1:A:254:THR:HG21	1.99	0.43
1:C:27:ILE:HD12	1:C:254:THR:HG21	2.00	0.43
1:B:87:GLY:C	1:B:89:GLY:H	2.27	0.43
1:B:166:ARG:O	1:B:170:GLN:HB2	2.19	0.43
1:A:329:ARG:O	1:A:333:LYS:HG3	2.19	0.43
1:A:171:THR:HG21	1:A:194:ILE:HG21	2.00	0.43
1:A:305:ARG:O	1:A:308:VAL:HG22	2.19	0.43
1:C:136:ARG:HB2	1:C:273:ILE:HG23	2.01	0.42
1:C:207:LEU:C	1:C:209:GLY:N	2.77	0.42
1:A:332:GLU:O	1:A:336:GLN:HG3	2.19	0.42
1:C:112:TRP:CD1	1:C:119:ALA:HB2	2.54	0.42
1:C:158:LEU:HB3	1:C:218:ASN:HD21	1.84	0.42
1:A:223:ASN:O	1:A:224:LYS:C	2.61	0.42
1:A:250:THR:O	1:A:265:PHE:HA	2.19	0.42
1:B:166:ARG:HB2	1:B:166:ARG:HE	1.74	0.42
1:B:211:ASN:HD22	1:B:212:THR:H	1.61	0.42
1:C:156:MET:HE2	1:C:156:MET:CA	2.49	0.42
1:B:223:ASN:O	1:B:224:LYS:C	2.62	0.42
1:C:288:GLN:HA	1:C:288:GLN:NE2	2.29	0.42
1:B:329:ARG:O	1:B:333:LYS:HG3	2.19	0.42
1:C:2:ALA:HA	5:C:1045:HOH:O	2.20	0.42
1:C:191:LEU:HD23	1:C:219:ASP:HB3	2.02	0.42
1:A:87:GLY:C	1:A:89:GLY:H	2.27	0.42
1:B:239:THR:HB	1:B:240:PRO:CD	2.50	0.42
1:C:112:TRP:CG	1:C:119:ALA:HB2	2.55	0.42
1:C:212:THR:HG23	1:C:213:PRO:N	2.33	0.42
1:C:240:PRO:C	1:C:242:GLU:N	2.77	0.42
1:A:20:GLU:O	1:A:61:LEU:HD13	2.20	0.42
1:B:12:ALA:HB2	1:B:111:ILE:HG13	2.01	0.41
1:B:161:ALA:HB1	1:B:169:ALA:HB3	2.02	0.41
1:B:170:GLN:HA	1:B:170:GLN:NE2	2.35	0.41
1:B:282:LEU:HD23	1:B:282:LEU:HA	1.94	0.41
1:A:136:ARG:HB2	1:A:273:ILE:HG23	2.01	0.41
1:A:157:ASP:C	1:A:159:GLY:N	2.78	0.41
1:B:156:MET:HE1	1:B:220:ILE:CG1	2.47	0.41
1:B:220:ILE:HA	1:B:230:ASN:HA	2.02	0.41
1:A:288:GLN:HE21	1:A:293:GLU:CG	2.34	0.41
1:B:199:PRO:HB3	1:B:214:VAL:HG11	2.03	0.41
1:A:190:ALA:C	1:A:191:LEU:HD12	2.46	0.41
1:A:236:PRO:O	1:A:237:GLU:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:HIS:HE1	1:C:160:HIS:CE1	2.38	0.41
1:B:192:MET:HE2	1:B:218:ASN:CB	2.45	0.41
1:B:225:VAL:HB	5:B:736:HOH:O	2.20	0.41
1:B:294:ASP:O	1:B:298:VAL:HG22	2.20	0.41
1:C:183:VAL:HG21	1:C:326:ARG:HD3	2.02	0.41
1:C:215:ASP:O	1:C:234:VAL:HA	2.21	0.41
1:C:220:ILE:HA	1:C:230:ASN:HA	2.03	0.41
1:C:263:TYR:CE1	1:C:289:ALA:HB1	2.55	0.41
1:A:151:LEU:HD23	1:A:151:LEU:HA	1.78	0.41
1:C:39:ASP:C	1:C:39:ASP:OD2	2.63	0.40
1:C:329:ARG:O	1:C:333:LYS:HG3	2.21	0.40
1:C:199:PRO:HG2	1:C:200:SER:N	2.35	0.40
1:C:341:ARG:NH1	5:C:1166:HOH:O	2.55	0.40
1:A:141:TYR:CG	1:A:142:GLY:N	2.87	0.40
1:B:156:MET:HE2	1:B:156:MET:CA	2.51	0.40
1:B:161:ALA:HA	1:B:165:HIS:CD2	2.57	0.40
1:B:304:ARG:HG3	1:B:304:ARG:NH1	2.37	0.40
1:B:305:ARG:O	1:B:308:VAL:HG22	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	339/349 (97%)	316 (93%)	17 (5%)	6 (2%)	6	4
1	B	339/349 (97%)	310 (91%)	25 (7%)	4 (1%)	10	8
1	C	323/349 (93%)	285 (88%)	24 (7%)	14 (4%)	2	1
All	All	1001/1047 (96%)	911 (91%)	66 (7%)	24 (2%)	4	3

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	VAL
1	C	158	LEU
1	C	208	ARG
1	C	225	VAL
1	A	211	ASN
1	B	166	ARG
1	B	225	VAL
1	C	159	GLY
1	C	160	HIS
1	C	237	GLU
1	A	224	LYS
1	A	237	GLU
1	B	224	LYS
1	C	180	GLU
1	C	198	THR
1	C	199	PRO
1	C	210	ALA
1	C	224	LYS
1	B	213	PRO
1	C	200	SER
1	C	203	MET
1	C	212	THR
1	A	196	GLY
1	A	212	THR

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	276/284 (97%)	271 (98%)	5 (2%)	51	68
1	B	276/284 (97%)	271 (98%)	5 (2%)	51	68
1	C	265/284 (93%)	258 (97%)	7 (3%)	40	55
All	All	817/852 (96%)	800 (98%)	17 (2%)	47	63

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	28	LEU
1	A	102	VAL
1	A	290	LEU
1	A	342	LEU
1	B	25	ARG
1	B	28	LEU
1	B	102	VAL
1	B	245	ILE
1	B	342	LEU
1	C	25	ARG
1	C	28	LEU
1	C	194	ILE
1	C	206	PHE
1	C	212	THR
1	C	215	ASP
1	C	291	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	63	ASN
1	A	143	HIS
1	A	168	ASN
1	A	170	GLN
1	A	211	ASN
1	A	223	ASN
1	A	288	GLN
1	B	63	ASN
1	B	143	HIS
1	B	154	ASN
1	B	162	GLN
1	B	170	GLN
1	B	211	ASN
1	B	223	ASN
1	B	288	GLN
1	C	63	ASN
1	C	83	HIS
1	C	143	HIS
1	C	160	HIS
1	C	243	GLN
1	C	286	GLN

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Mol	Chain	Res	Type
1	C	288	GLN

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](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	FES	C	501	1	0,4,4	-	-	-		
4	D3M	B	601	-	13,13,13	1.80	3 (23%)	18,18,18	1.18	2 (11%)
2	FES	A	501	1	0,4,4	-	-	-		
4	D3M	A	601	-	13,13,13	1.67	1 (7%)	18,18,18	1.13	2 (11%)
2	FES	B	501	1	0,4,4	-	-	-		

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	FES	C	501	1	-	-	0/1/1/1
4	D3M	B	601	-	-	0/6/6/6	0/1/1/1
2	FES	A	501	1	-	-	0/1/1/1
4	D3M	A	601	-	-	0/6/6/6	0/1/1/1
2	FES	B	501	1	-	-	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	601	D3M	C1-C2	3.71	1.45	1.39
4	A	601	D3M	C1-C2	3.50	1.44	1.39
4	B	601	D3M	O3-C6	2.33	1.42	1.38
4	B	601	D3M	C1-C6	2.13	1.45	1.40

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	D3M	C6-C5-CL1	3.00	122.36	118.40
4	A	601	D3M	C6-C5-CL1	2.72	121.98	118.40
4	B	601	D3M	C1-C2-CL2	2.29	122.56	119.62
4	A	601	D3M	C1-C2-CL2	2.12	122.35	119.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	FES	1	0
2	A	501	FES	1	0
2	B	501	FES	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

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	341/349 (97%)	0.13	5 (1%) 72 69	26, 43, 68, 89	0
1	B	341/349 (97%)	0.31	21 (6%) 26 23	29, 45, 85, 90	0
1	C	327/349 (93%)	0.34	29 (8%) 15 13	27, 42, 85, 90	0
All	All	1009/1047 (96%)	0.26	55 (5%) 30 27	26, 43, 82, 90	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	202	LEU	5.3
1	B	342	LEU	4.0
1	C	199	PRO	3.8
1	C	201	VAL	3.7
1	C	198	THR	3.7
1	B	210	ALA	3.6
1	B	240	PRO	3.5
1	C	339	ALA	3.4
1	C	2	ALA	3.3
1	C	200	SER	3.3
1	B	212	THR	3.3
1	C	212	THR	3.3
1	C	206	PHE	3.2
1	C	177	LEU	3.2
1	B	63	ASN	3.1
1	C	196	GLY	3.1
1	C	211	ASN	3.1
1	B	172	ASP	3.0
1	C	160	HIS	3.0
1	C	197	GLY	2.9
1	C	203	MET	2.9
1	C	158	LEU	2.9
1	A	209	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	195	PRO	2.8
1	A	225	VAL	2.8
1	B	209	GLY	2.7
1	A	2	ALA	2.7
1	C	276	PRO	2.7
1	C	207	LEU	2.7
1	B	2	ALA	2.7
1	B	273	ILE	2.7
1	B	239	THR	2.6
1	B	211	ASN	2.6
1	C	278	MET	2.5
1	B	195	PRO	2.5
1	B	213	PRO	2.5
1	B	271	PHE	2.5
1	A	3	THR	2.5
1	A	62	VAL	2.4
1	C	213	PRO	2.4
1	C	238	GLY	2.4
1	B	236	PRO	2.4
1	B	194	ILE	2.4
1	C	209	GLY	2.3
1	C	162	GLN	2.3
1	B	225	VAL	2.3
1	C	337	LEU	2.3
1	B	186	GLY	2.2
1	B	214	VAL	2.2
1	C	225	VAL	2.1
1	B	159	GLY	2.1
1	C	215	ASP	2.1
1	C	161	ALA	2.0
1	C	204	ALA	2.0
1	B	198	THR	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 [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
4	D3M	B	601	13/13	0.86	0.10	45,49,50,51	0
4	D3M	A	601	13/13	0.91	0.09	29,37,41,41	0
3	FE	B	502	1/1	0.95	0.06	54,54,54,54	0
2	FES	A	501	4/4	0.98	0.04	37,40,40,41	0
2	FES	B	501	4/4	0.99	0.04	29,30,30,31	0
2	FES	C	501	4/4	0.99	0.02	28,29,29,31	0
3	FE	A	502	1/1	0.99	0.01	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.