



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:00 AM UTC

PDB ID : 1M1T / pdb_00001m1t
Title : Biosynthetic thiolase, Q64A mutant
Authors : Kursula, P.; Ojala, J.; Lambeir, A.-M.; Wierenga, R.K.
Deposited on : 2002-06-20
Resolution : 1.94 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

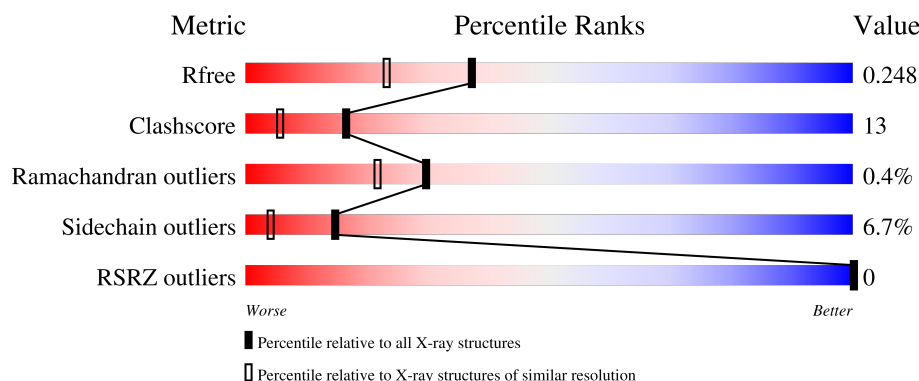
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	392	 77% 21% .
1	B	392	 74% 22% . .
1	C	392	 78% 21% .
1	D	392	 75% 22% .

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
3	GOL	B	7393	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12545 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.

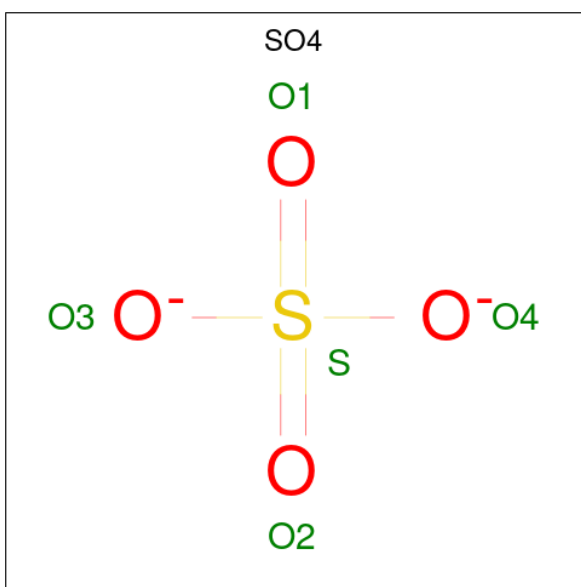
- Molecule 1 is a protein called Acetyl-CoA acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	1	0
			2831	1757	511	542	21			
1	B	392	Total	C	N	O	S	0	1	0
			2831	1757	511	542	21			
1	C	392	Total	C	N	O	S	0	1	0
			2831	1757	511	542	21			
1	D	392	Total	C	N	O	S	0	1	0
			2831	1757	511	542	21			

There are 12 discrepancies between the modelled and reference sequences:

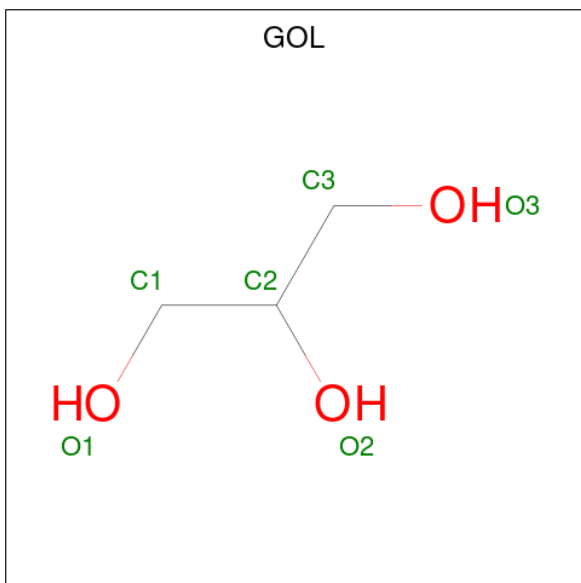
Chain	Residue	Modelled	Actual	Comment	Reference
A	10	ALA	-	insertion	UNP P07097
A	64	ALA	GLN	engineered mutation	UNP P07097
A	129	ARG	ALA	conflict	UNP P07097
B	10	ALA	-	insertion	UNP P07097
B	64	ALA	GLN	engineered mutation	UNP P07097
B	129	ARG	ALA	conflict	UNP P07097
C	10	ALA	-	insertion	UNP P07097
C	64	ALA	GLN	engineered mutation	UNP P07097
C	129	ARG	ALA	conflict	UNP P07097
D	10	ALA	-	insertion	UNP P07097
D	64	ALA	GLN	engineered mutation	UNP P07097
D	129	ARG	ALA	conflict	UNP P07097

- 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		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

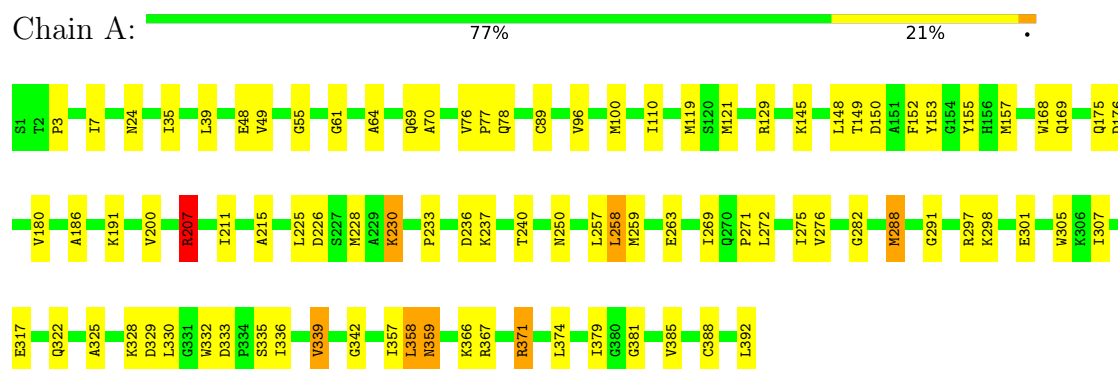
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	424	Total	O	0	0
			424	424		
4	B	440	Total	O	0	0
			440	440		
4	C	151	Total	O	0	0
			151	151		
4	D	162	Total	O	0	0
			162	162		

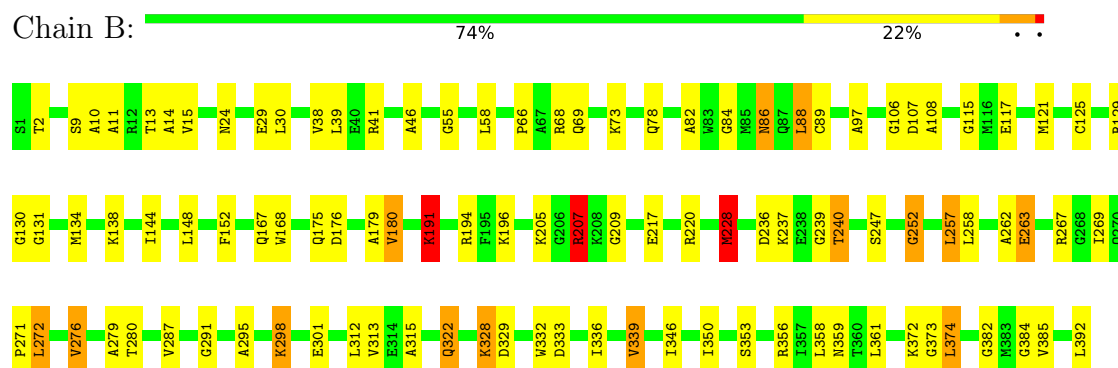
3 Residue-property plots

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.

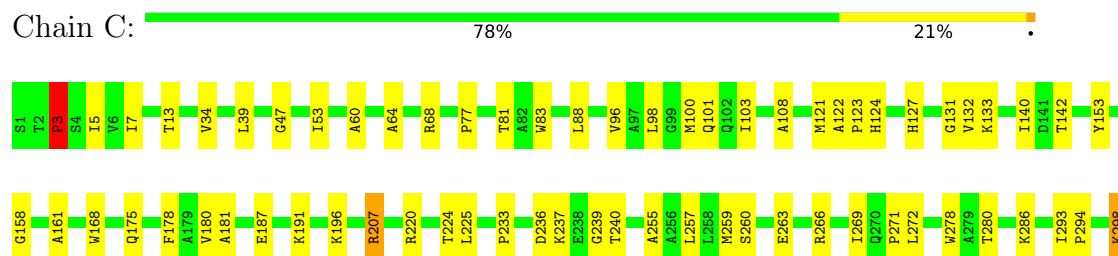
• Molecule 1: Acetyl-CoA acetyltransferase

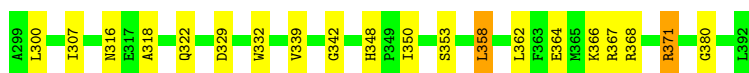


• Molecule 1: Acetyl-CoA acetyltransferase



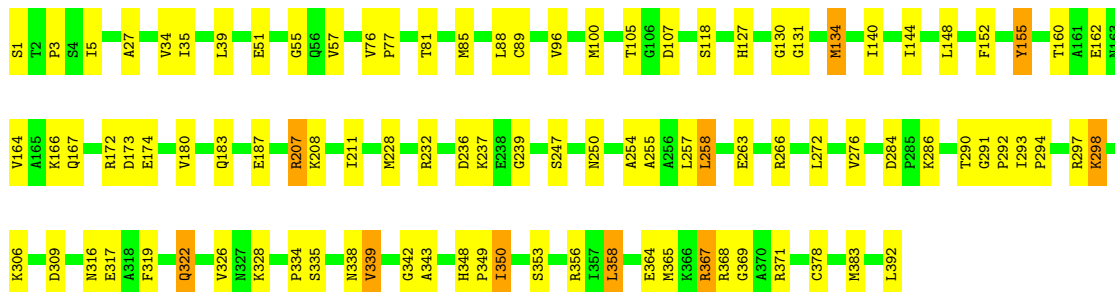
• Molecule 1: Acetyl-CoA acetyltransferase





• Molecule 1: Acetyl-CoA acetyltransferase

Chain D: 75% 22% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.47Å 78.95Å 148.29Å 90.00° 92.55° 90.00°	Depositor
Resolution (Å)	20.00 – 1.94 20.00 – 1.94	Depositor EDS
% Data completeness (in resolution range)	97.5 (20.00-1.94) 85.4 (20.00-1.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 1.94Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.199 , 0.245 0.208 , 0.248	Depositor DCC
R_{free} test set	6356 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.828	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.137 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12545	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.45	11/2877 (0.4%)	1.31	8/3886 (0.2%)
1	B	1.50	28/2877 (1.0%)	1.35	11/3886 (0.3%)
1	C	0.85	0/2877	1.11	4/3886 (0.1%)
1	D	0.82	0/2877	1.09	2/3886 (0.1%)
All	All	1.20	39/11508 (0.3%)	1.22	25/15544 (0.2%)

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	315	ALA	CA-CB	7.50	1.63	1.53
1	B	11	ALA	CA-CB	7.38	1.65	1.53
1	B	97	ALA	CA-CB	6.87	1.64	1.53
1	B	313	VAL	C-O	-6.81	1.17	1.23
1	A	215	ALA	CA-C	6.36	1.60	1.52
1	A	70	ALA	CA-CB	6.25	1.63	1.53
1	B	382	GLY	C-O	6.20	1.32	1.23
1	B	262	ALA	CA-CB	-6.19	1.43	1.53
1	B	46	ALA	C-O	6.17	1.31	1.24
1	B	68	ARG	CA-C	-6.06	1.45	1.52
1	B	14	ALA	CA-CB	-6.03	1.43	1.53
1	A	357	ILE	CA-CB	5.99	1.62	1.54
1	B	179	ALA	CA-CB	5.99	1.62	1.53
1	B	38	VAL	CA-CB	5.78	1.61	1.54
1	B	228	MET	SD-CE	-5.76	1.65	1.79
1	A	49	VAL	C-O	-5.66	1.16	1.24
1	B	372	LYS	CA-C	5.63	1.59	1.52
1	B	108	ALA	CA-CB	5.63	1.64	1.54
1	B	295	ALA	CA-CB	-5.60	1.44	1.53
1	B	276	VAL	C-O	-5.54	1.17	1.24
1	B	82	ALA	CA-CB	-5.53	1.44	1.53
1	A	35	ILE	CA-C	5.48	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	ALA	CA-CB	-5.47	1.44	1.53
1	B	46	ALA	CA-CB	5.34	1.62	1.53
1	B	385	VAL	CA-CB	-5.33	1.47	1.54
1	A	211	ILE	CA-CB	5.32	1.61	1.54
1	B	117	GLU	CA-C	-5.31	1.45	1.52
1	A	70	ALA	CA-C	5.29	1.59	1.52
1	B	374	LEU	C-O	-5.24	1.17	1.23
1	A	335	SER	C-O	-5.22	1.17	1.24
1	B	84	GLY	CA-C	-5.20	1.47	1.52
1	B	191	LYS	C-O	-5.18	1.18	1.24
1	B	194	ARG	C-O	-5.18	1.17	1.24
1	B	115	GLY	CA-C	-5.11	1.46	1.51
1	A	288	MET	C-O	-5.10	1.17	1.24
1	A	119	MET	N-CA	-5.08	1.40	1.46
1	B	58	LEU	C-O	-5.08	1.18	1.23
1	B	88	LEU	C-O	-5.06	1.19	1.24
1	B	10	ALA	CA-C	-5.04	1.46	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	THR	N-CA-C	-7.37	92.58	108.64
1	A	371	ARG	N-CA-C	6.58	118.24	111.14
1	A	55	GLY	N-CA-C	-6.32	102.94	111.54
1	A	359	ASN	N-CA-C	-6.25	104.39	111.14
1	C	3	PRO	N-CA-C	6.24	125.33	112.47
1	B	252	GLY	N-CA-C	5.84	119.28	110.18
1	D	140	ILE	N-CA-C	5.70	116.52	108.36
1	B	107	ASP	N-CA-C	5.65	119.89	112.89
1	D	250	ASN	N-CA-C	5.64	116.83	108.60
1	C	13	THR	N-CA-C	-5.49	103.24	110.43
1	B	207	ARG	N-CA-C	-5.41	104.72	112.13
1	B	55	GLY	N-CA-C	-5.40	104.20	111.54
1	A	339	VAL	N-CA-CB	-5.35	101.08	110.77
1	B	339	VAL	N-CA-CB	-5.35	101.51	110.65
1	A	339	VAL	CG1-CB-CG2	5.32	122.50	110.80
1	B	86	ASN	N-CA-C	5.31	117.70	107.44
1	B	130	GLY	N-CA-C	-5.28	107.78	114.16
1	B	180	VAL	CB-CA-C	-5.26	105.03	112.14
1	B	373	GLY	N-CA-C	5.14	120.51	111.57
1	C	178	PHE	N-CA-C	-5.13	105.60	111.14
1	B	129	ARG	N-CA-C	5.12	116.86	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	250	ASN	N-CA-C	5.12	116.58	108.96
1	A	207	ARG	N-CA-C	-5.07	104.74	112.04
1	A	291	GLY	N-CA-C	5.06	122.66	112.34
1	C	131	GLY	N-CA-C	-5.03	106.24	112.33

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2831	0	2839	72	0
1	B	2831	0	2839	77	0
1	C	2831	0	2839	66	0
1	D	2831	0	2839	90	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	12	0	16	3	0
3	B	12	0	16	7	0
4	A	424	0	0	47	1
4	B	440	0	0	40	2
4	C	151	0	0	23	0
4	D	162	0	0	49	0
All	All	12545	0	11388	294	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (294) 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:C:60:ALA:HB3	4:C:492:HOH:O	1.60	0.98
1:A:258:LEU:HG	4:A:6756:HOH:O	1.65	0.96
1:D:76:VAL:HG23	4:D:534:HOH:O	1.63	0.95
1:D:27:ALA:HB2	4:D:443:HOH:O	1.67	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:PHE:C	4:A:6720:HOH:O	2.10	0.94
1:A:100:MET:HG3	4:A:6747:HOH:O	1.67	0.93
1:B:252:GLY:HA3	4:B:7826:HOH:O	1.67	0.92
1:C:207:ARG:HA	4:C:539:HOH:O	1.69	0.92
1:C:133:LYS:HB2	4:C:528:HOH:O	1.71	0.90
1:A:157:MET:HG3	4:A:6719:HOH:O	1.72	0.89
1:A:61:GLY:HA2	4:A:6758:HOH:O	1.72	0.88
1:D:356:ARG:HA	4:D:542:HOH:O	1.73	0.87
1:D:392:LEU:HD21	4:D:448:HOH:O	1.75	0.87
1:B:247[A]:SER:OG	3:B:7393:GOL:H31	1.77	0.85
1:D:208:LYS:HD3	4:D:540:HOH:O	1.78	0.82
1:D:338:ASN:ND2	4:D:500:HOH:O	2.12	0.82
1:A:48:GLU:CD	4:A:6767:HOH:O	2.25	0.80
1:D:290:THR:HA	4:D:511:HOH:O	1.82	0.79
1:C:181:ALA:HB3	4:C:543:HOH:O	1.84	0.78
1:B:88:LEU:HD13	4:B:7746:HOH:O	1.83	0.77
1:C:187:GLU:HG2	4:C:533:HOH:O	1.85	0.77
1:B:346:ILE:HG12	4:B:7776:HOH:O	1.86	0.74
1:C:5:ILE:HG12	4:C:523:HOH:O	1.86	0.74
1:B:9:SER:HB2	4:B:7545:HOH:O	1.86	0.74
1:D:254:ALA:HB1	4:D:542:HOH:O	1.87	0.74
1:B:207:ARG:HD2	4:B:7813:HOH:O	1.88	0.74
4:A:6817:HOH:O	1:C:133:LYS:CD	2.35	0.74
1:B:138:LYS:NZ	4:B:7705:HOH:O	2.19	0.74
1:A:61:GLY:CA	4:A:6758:HOH:O	2.32	0.73
4:A:6758:HOH:O	1:B:125:CYS:SG	2.47	0.73
1:B:374:LEU:HD23	1:B:374:LEU:C	2.13	0.72
1:A:258:LEU:HG	4:A:6436:HOH:O	1.88	0.72
1:A:207:ARG:HG2	1:A:207:ARG:HH11	1.55	0.71
1:B:152:PHE:CE2	4:B:7757:HOH:O	2.43	0.71
1:B:13:THR:HB	4:B:7701:HOH:O	1.93	0.69
1:D:167:GLN:HG2	4:D:397:HOH:O	1.92	0.69
1:D:207:ARG:CZ	4:D:464:HOH:O	2.39	0.69
1:A:288:MET:SD	4:A:6710:HOH:O	2.50	0.69
1:B:15:VAL:HA	4:B:7826:HOH:O	1.92	0.69
3:B:7393:GOL:H32	4:B:7817:HOH:O	1.92	0.69
1:C:3:PRO:HG2	4:C:508:HOH:O	1.93	0.69
1:A:48:GLU:OE2	4:A:6767:HOH:O	2.10	0.68
1:D:369:GLY:HA3	4:D:515:HOH:O	1.92	0.68
1:B:257:LEU:HD23	1:B:258:LEU:N	2.09	0.68
1:D:152:PHE:CE2	4:D:503:HOH:O	2.47	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:PHE:CD2	4:B:7750:HOH:O	2.46	0.68
1:B:175:GLN:HE22	1:B:240:THR:CG2	2.07	0.68
1:A:200:VAL:HG23	3:A:6394:GOL:H2	1.76	0.67
1:C:196:LYS:NZ	4:C:426:HOH:O	2.27	0.67
1:D:127:HIS:CG	4:D:539:HOH:O	2.47	0.66
1:A:152:PHE:HD2	4:A:6743:HOH:O	1.77	0.66
1:A:329:ASP:HA	4:A:6784:HOH:O	1.94	0.66
1:C:175:GLN:HE22	1:C:240:THR:CG2	2.09	0.66
1:A:175:GLN:HE22	1:A:240:THR:CG2	2.08	0.66
1:A:330:LEU:HD22	4:A:6815:HOH:O	1.95	0.66
1:D:35:ILE:HG22	4:D:534:HOH:O	1.95	0.65
1:B:13:THR:CB	4:B:7701:HOH:O	2.45	0.65
1:D:334:PRO:O	4:D:480:HOH:O	2.14	0.65
1:C:367:ARG:HD3	4:C:531:HOH:O	1.95	0.65
1:A:225:LEU:HD21	4:A:6722:HOH:O	1.95	0.65
4:A:6817:HOH:O	1:C:133:LYS:HD2	1.96	0.65
1:B:13:THR:CG2	4:B:7701:HOH:O	2.45	0.65
1:D:284:ASP:HB2	4:D:547:HOH:O	1.97	0.65
1:B:30:LEU:HD22	4:B:7701:HOH:O	1.97	0.64
1:D:319:PHE:HB2	4:D:502:HOH:O	1.97	0.64
1:A:301:GLU:HG2	4:A:6790:HOH:O	1.96	0.64
1:B:29:GLU:OE1	4:B:7670:HOH:O	2.15	0.64
1:C:101:GLN:HG2	1:D:105:THR:HG21	1.80	0.63
1:D:297:ARG:NE	4:D:486:HOH:O	2.31	0.63
1:A:100:MET:HE2	4:A:6747:HOH:O	1.98	0.63
1:C:263:GLU:OE1	1:C:266:ARG:NH1	2.31	0.63
1:B:168:TRP:CH2	1:B:329:ASP:HB2	2.34	0.63
1:D:207:ARG:N	1:D:207:ARG:HD3	2.14	0.62
1:C:207:ARG:HG2	1:C:207:ARG:HH11	1.64	0.62
1:A:307:ILE:HD11	4:A:6815:HOH:O	1.99	0.62
3:B:7394:GOL:H31	4:B:7796:HOH:O	2.00	0.62
1:D:118:SER:HB3	4:D:443:HOH:O	1.98	0.62
1:A:169:GLN:HG3	4:A:6537:HOH:O	1.99	0.62
1:B:209:GLY:HA2	4:B:7829:HOH:O	1.99	0.62
1:C:133:LYS:CB	4:C:528:HOH:O	2.38	0.61
1:D:236:ASP:HB2	4:D:538:HOH:O	2.00	0.61
1:B:301:GLU:HG2	4:B:7735:HOH:O	1.99	0.61
1:B:175:GLN:HE22	1:B:240:THR:HG23	1.65	0.61
1:A:152:PHE:CE1	1:B:69:GLN:HA	2.34	0.61
1:A:366:LYS:HE2	4:A:6818:HOH:O	2.00	0.61
1:A:275:ILE:HG13	4:A:6809:HOH:O	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ARG:CD	4:B:7813:HOH:O	2.46	0.60
1:B:392:LEU:HD21	4:B:7665:HOH:O	2.01	0.60
1:B:191:LYS:NZ	4:B:7649:HOH:O	2.33	0.60
4:A:6817:HOH:O	1:C:133:LYS:HD3	1.98	0.60
1:C:207:ARG:H	1:C:207:ARG:HD3	1.66	0.60
1:B:205:LYS:HE2	4:B:7821:HOH:O	2.01	0.60
1:B:167:GLN:HA	4:B:7791:HOH:O	2.02	0.59
1:D:183:GLN:HG3	4:D:545:HOH:O	2.02	0.59
4:A:6741:HOH:O	1:D:134:MET:CE	2.50	0.59
1:C:53:ILE:HG12	1:C:83:TRP:CE2	2.38	0.59
1:C:98:LEU:CD2	4:D:465:HOH:O	2.50	0.58
1:B:13:THR:HG21	4:B:7701:HOH:O	2.01	0.58
1:D:263:GLU:OE1	1:D:266:ARG:NH1	2.36	0.58
1:C:142:THR:HG21	4:C:492:HOH:O	2.03	0.58
1:B:279:ALA:HB1	1:B:298:LYS:HG3	1.85	0.58
1:C:371:ARG:HD2	4:C:444:HOH:O	2.03	0.58
1:D:77:PRO:HD2	4:D:541:HOH:O	2.04	0.58
1:D:335:SER:HA	4:D:482:HOH:O	2.03	0.58
1:B:247[A]:SER:OG	3:B:7393:GOL:C3	2.50	0.58
1:D:180:VAL:HA	4:D:545:HOH:O	2.04	0.58
1:A:150:ASP:OD1	4:A:6743:HOH:O	2.18	0.57
1:C:5:ILE:HG13	1:C:100:MET:HG2	1.85	0.57
1:A:152:PHE:CE2	4:A:6667:HOH:O	2.52	0.57
1:A:175:GLN:HE22	1:A:240:THR:HG23	1.68	0.57
1:B:346:ILE:HG23	4:B:7776:HOH:O	2.03	0.57
1:C:133:LYS:NZ	4:C:432:HOH:O	2.37	0.57
1:D:130:GLY:N	4:D:452:HOH:O	2.38	0.57
1:B:228:MET:SD	4:B:7658:HOH:O	2.58	0.56
1:C:207:ARG:HD3	1:C:207:ARG:N	2.20	0.56
1:A:258:LEU:CG	4:A:6756:HOH:O	2.38	0.56
1:B:207:ARG:HD3	1:B:207:ARG:N	2.20	0.56
1:A:152:PHE:HE1	1:B:69:GLN:HA	1.71	0.56
1:A:257:LEU:C	1:A:258:LEU:HD22	2.30	0.56
1:D:35:ILE:CG2	4:D:534:HOH:O	2.52	0.56
1:A:230:LYS:NZ	4:A:6705:HOH:O	2.36	0.56
1:D:55:GLY:HA2	1:D:85:MET:O	2.06	0.56
1:B:205:LYS:CE	4:B:7821:HOH:O	2.54	0.55
1:D:328:LYS:HE2	4:D:491:HOH:O	2.05	0.55
1:D:232:ARG:CZ	4:D:523:HOH:O	2.54	0.55
1:D:207:ARG:HG2	1:D:207:ARG:HH11	1.71	0.55
1:C:60:ALA:CB	4:C:492:HOH:O	2.33	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:162:GLU:HG3	4:D:538:HOH:O	2.05	0.55
1:D:297:ARG:CZ	4:D:486:HOH:O	2.55	0.55
1:A:153:TYR:HD1	4:A:6682:HOH:O	1.91	0.54
1:C:237:LYS:HD3	4:C:522:HOH:O	2.08	0.54
1:C:121:MET:CE	4:D:539:HOH:O	2.56	0.54
1:C:339:VAL:HG11	1:C:368:ARG:NH2	2.22	0.53
1:A:297:ARG:HD3	4:A:6623:HOH:O	2.09	0.53
1:B:322:GLN:HB3	4:B:7814:HOH:O	2.09	0.53
1:C:269:ILE:O	1:C:271:PRO:HD3	2.08	0.53
1:D:257:LEU:HD23	1:D:257:LEU:C	2.33	0.53
1:C:175:GLN:HE22	1:C:240:THR:HG21	1.73	0.53
1:A:78:GLN:NE2	4:A:6744:HOH:O	2.42	0.53
1:A:3:PRO:HD3	4:A:6804:HOH:O	2.09	0.52
1:D:316:ASN:HA	4:D:529:HOH:O	2.09	0.52
1:D:166:LYS:NZ	4:D:460:HOH:O	2.31	0.52
1:D:172:ARG:HB2	4:D:492:HOH:O	2.09	0.52
1:D:211:ILE:HB	4:D:554:HOH:O	2.10	0.52
1:D:257:LEU:HD23	1:D:258:LEU:N	2.24	0.52
1:A:333:ASP:O	1:A:336:ILE:HG12	2.10	0.52
1:D:237:LYS:HA	4:D:444:HOH:O	2.09	0.52
1:A:168:TRP:CH2	1:A:329:ASP:HB2	2.45	0.52
1:B:168:TRP:HH2	1:B:329:ASP:HB2	1.75	0.52
1:A:258:LEU:HB2	4:A:6756:HOH:O	2.10	0.51
1:C:280:THR:HG23	1:D:81:THR:HG21	1.92	0.51
1:A:157:MET:CG	4:A:6719:HOH:O	2.45	0.51
1:A:258:LEU:CB	4:A:6756:HOH:O	2.57	0.51
1:C:122:ALA:HA	4:C:434:HOH:O	2.09	0.51
1:B:237:LYS:N	1:B:237:LYS:HD2	2.26	0.51
1:A:258:LEU:HD22	1:A:258:LEU:N	2.24	0.51
1:D:236:ASP:HB3	1:D:239:GLY:HA3	1.92	0.51
1:D:339:VAL:HG11	1:D:368:ARG:NH2	2.26	0.51
1:D:364:GLU:OE1	1:D:367:ARG:NH1	2.40	0.51
1:A:176:ASP:O	1:A:180:VAL:HG23	2.11	0.50
1:A:207:ARG:HH11	1:A:207:ARG:CG	2.21	0.50
1:C:123:PRO:HG2	4:C:492:HOH:O	2.11	0.50
1:A:371:ARG:HD2	4:A:6788:HOH:O	2.11	0.50
1:C:34:VAL:HG12	1:C:255:ALA:HB3	1.93	0.50
1:D:27:ALA:CB	4:D:443:HOH:O	2.39	0.50
1:B:228:MET:CE	4:B:7658:HOH:O	2.59	0.50
1:B:328:LYS:NZ	4:B:7745:HOH:O	2.05	0.50
1:B:217:GLU:HB3	4:B:7508:HOH:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:LEU:HD23	1:B:361:LEU:HD12	1.94	0.50
1:B:15:VAL:HG22	4:B:7826:HOH:O	2.11	0.49
1:A:69:GLN:HA	1:B:152:PHE:CE1	2.48	0.49
1:B:207:ARG:HB2	4:B:7783:HOH:O	2.12	0.49
1:D:57:VAL:HG21	1:D:350:ILE:HG22	1.95	0.49
1:A:175:GLN:HE22	1:A:240:THR:HG21	1.77	0.49
1:A:297:ARG:HG2	1:A:297:ARG:HH11	1.78	0.49
1:C:278:TRP:NE1	1:D:107:ASP:OD2	2.43	0.49
1:D:164:VAL:HG21	4:D:469:HOH:O	2.12	0.49
1:B:196:LYS:NZ	4:B:7555:HOH:O	2.46	0.49
1:A:305:TRP:CZ3	1:A:388:CYS:HB3	2.48	0.49
1:B:86:ASN:OD1	1:B:86:ASN:C	2.55	0.49
1:B:175:GLN:HE22	1:B:240:THR:HG21	1.76	0.49
1:B:333:ASP:O	1:B:336:ILE:HG12	2.12	0.48
1:C:236:ASP:HB3	1:C:239:GLY:HA3	1.96	0.48
1:D:127:HIS:CB	4:D:539:HOH:O	2.61	0.48
1:D:291:GLY:N	1:D:292:PRO:CD	2.76	0.48
1:C:364:GLU:O	1:C:368:ARG:HG2	2.13	0.48
1:B:269:ILE:O	1:B:271:PRO:HD3	2.12	0.48
1:D:144:ILE:HD13	1:D:148:LEU:HD12	1.94	0.48
1:B:176:ASP:O	1:B:180:VAL:HG23	2.13	0.48
1:C:168:TRP:CH2	1:C:329:ASP:HB2	2.49	0.48
1:D:51:GLU:HA	1:D:81:THR:O	2.13	0.48
1:B:263:GLU:CD	1:B:267:ARG:HE	2.21	0.48
1:C:124:HIS:HA	1:C:140:ILE:O	2.14	0.48
1:A:145:LYS:NZ	4:A:6664:HOH:O	2.41	0.48
1:C:318:ALA:CB	4:C:458:HOH:O	2.61	0.47
1:C:318:ALA:HB1	4:C:458:HOH:O	2.13	0.47
1:A:233:PRO:HB2	1:A:236:ASP:O	2.14	0.47
1:C:68:ARG:HG3	1:D:152:PHE:HZ	1.80	0.47
1:C:88:LEU:HD12	1:C:380:GLY:O	2.14	0.47
1:A:153:TYR:HB3	1:A:155:TYR:CE2	2.49	0.47
1:D:349:PRO:O	1:D:350:ILE:C	2.57	0.47
1:A:325:ALA:HB1	4:A:6771:HOH:O	2.14	0.47
1:B:41:ARG:HE	3:B:7394:GOL:H12	1.79	0.47
1:A:282:GLY:HA3	1:B:78:GLN:O	2.15	0.47
1:D:317:GLU:CD	1:D:342:GLY:HA3	2.40	0.47
1:A:152:PHE:CD2	4:A:6743:HOH:O	2.55	0.46
1:D:174:GLU:OE2	1:D:328:LYS:NZ	2.45	0.46
1:D:228:MET:CE	4:D:536:HOH:O	2.63	0.46
1:B:152:PHE:CG	4:B:7750:HOH:O	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:ILE:HA	1:A:258:LEU:HD13	1.96	0.46
1:D:292:PRO:HD3	1:D:378:CYS:HB3	1.97	0.46
1:C:81:THR:HG22	1:D:383:MET:HG2	1.97	0.46
1:C:257:LEU:HD22	1:C:259:MET:HE3	1.97	0.46
1:A:381:GLY:HA2	4:A:6743:HOH:O	2.15	0.45
1:D:5:ILE:HG13	1:D:100:MET:HG2	1.97	0.45
1:B:24:ASN:HA	1:B:121:MET:SD	2.57	0.45
1:D:34:VAL:CG1	1:D:255:ALA:HB3	2.45	0.45
1:D:160:THR:HB	1:D:286:LYS:O	2.17	0.45
1:A:76:VAL:HG13	1:A:77:PRO:HD2	1.98	0.44
1:C:7:ILE:HD13	1:C:362:LEU:HD11	1.99	0.44
1:A:168:TRP:HZ3	4:A:6784:HOH:O	1.99	0.44
1:A:269:ILE:O	1:A:271:PRO:HD3	2.17	0.44
1:A:148:LEU:O	1:A:157:MET:HG2	2.17	0.44
1:C:127:HIS:CE1	4:C:416:HOH:O	2.70	0.44
1:D:326:VAL:HG23	4:D:511:HOH:O	2.16	0.44
1:B:356:ARG:C	1:B:356:ARG:HD2	2.43	0.44
1:D:57:VAL:HG21	1:D:350:ILE:CG2	2.47	0.44
1:B:106:GLY:HA2	4:B:7578:HOH:O	2.17	0.44
1:B:144:ILE:HD13	1:B:148:LEU:HD12	1.98	0.44
3:A:6394:GOL:O3	4:A:6529:HOH:O	2.21	0.43
1:C:293:ILE:HB	1:C:294:PRO:CD	2.48	0.43
1:A:24:ASN:HA	1:A:121:MET:SD	2.57	0.43
1:A:379:ILE:HD11	1:A:385:VAL:HG12	2.00	0.43
1:D:322:GLN:O	1:D:326:VAL:HG23	2.18	0.43
1:B:228:MET:HE3	4:B:7658:HOH:O	2.18	0.43
1:C:64:ALA:HB1	1:D:88:LEU:HD11	2.01	0.43
1:C:233:PRO:HB2	1:C:236:ASP:O	2.17	0.43
1:D:207:ARG:HD3	1:D:207:ARG:H	1.83	0.43
1:D:306:LYS:O	1:D:309:ASP:HB2	2.19	0.43
1:D:173:ASP:HB3	4:D:528:HOH:O	2.18	0.43
1:A:374:LEU:C	1:A:374:LEU:HD23	2.44	0.43
1:B:131:GLY:HA2	1:D:131:GLY:HA2	2.01	0.43
1:B:228:MET:N	1:B:228:MET:HE2	2.34	0.43
1:D:293:ILE:HB	1:D:294:PRO:CD	2.49	0.43
1:B:41:ARG:HE	3:B:7394:GOL:C1	2.31	0.43
1:D:247[B]:SER:OG	1:D:348:HIS:HB2	2.19	0.43
1:B:257:LEU:C	1:B:257:LEU:CD2	2.92	0.43
1:B:287:VAL:HG21	4:B:7742:HOH:O	2.18	0.43
1:D:257:LEU:C	1:D:257:LEU:CD2	2.91	0.43
1:C:342:GLY:HA2	4:C:504:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:232:ARG:NH1	4:D:523:HOH:O	2.52	0.42
1:D:254:ALA:CB	4:D:542:HOH:O	2.56	0.42
1:B:280:THR:HA	1:B:384:GLY:O	2.19	0.42
1:C:207:ARG:HG2	1:C:207:ARG:NH1	2.32	0.42
1:D:228:MET:HE2	4:D:536:HOH:O	2.19	0.42
1:C:158:GLY:O	1:C:161:ALA:HB3	2.19	0.42
1:C:180:VAL:HG21	1:C:225:LEU:HA	2.01	0.42
1:B:236:ASP:HB3	1:B:239:GLY:HA3	2.01	0.42
1:A:330:LEU:HD23	1:A:330:LEU:HA	1.79	0.42
3:A:6394:GOL:H12	4:A:6734:HOH:O	2.18	0.42
1:C:298:LYS:HD3	1:C:298:LYS:C	2.44	0.42
1:B:257:LEU:HD23	1:B:257:LEU:C	2.44	0.42
1:B:73:LYS:CE	4:B:7806:HOH:O	2.67	0.42
1:C:316:ASN:HB3	4:C:517:HOH:O	2.19	0.42
1:D:364:GLU:O	1:D:368:ARG:HG2	2.19	0.42
1:C:121:MET:HA	1:D:127:HIS:CD2	2.55	0.42
1:A:64:ALA:HB1	1:B:88:LEU:HD11	2.00	0.42
1:D:155:TYR:HE1	1:D:160:THR:HG22	1.85	0.42
1:B:41:ARG:HB2	4:B:7545:HOH:O	2.20	0.41
1:D:152:PHE:CZ	4:D:503:HOH:O	2.72	0.41
1:B:207:ARG:HG2	1:B:207:ARG:HH11	1.85	0.41
1:C:298:LYS:HG3	4:C:521:HOH:O	2.20	0.41
1:C:348:HIS:CE1	1:C:353:SER:HG	2.35	0.41
1:D:96:VAL:HG21	1:D:358:LEU:HD12	2.02	0.41
1:A:317:GLU:CD	1:A:342:GLY:HA3	2.45	0.41
1:B:9:SER:HA	1:B:272:LEU:HD22	2.02	0.41
1:A:7:ILE:HD12	4:A:6809:HOH:O	2.20	0.41
1:D:317:GLU:O	1:D:343:ALA:HB3	2.21	0.41
1:A:275:ILE:CG2	4:A:6747:HOH:O	2.69	0.41
1:B:247[A]:SER:HG	3:B:7393:GOL:H31	1.80	0.41
1:C:96:VAL:HG21	1:C:358:LEU:HD12	2.02	0.41
1:C:153:TYR:CE2	1:C:286:LYS:HD3	2.56	0.41
1:D:207:ARG:HG2	1:D:207:ARG:NH1	2.35	0.41
1:A:96:VAL:HG21	1:A:358:LEU:HD12	2.02	0.41
1:A:257:LEU:CD2	1:A:259:MET:HE3	2.51	0.41
1:C:47:GLY:HA2	1:C:77:PRO:HG3	2.03	0.41
1:D:284:ASP:N	4:D:547:HOH:O	2.52	0.41
1:A:129:ARG:HA	1:C:132:VAL:O	2.21	0.40
1:A:69:GLN:HA	1:B:152:PHE:HE1	1.85	0.40
4:A:6741:HOH:O	1:D:134:MET:HE1	2.20	0.40
1:C:103:ILE:HA	1:C:108:ALA:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:294:PRO:O	1:D:298:LYS:HB2	2.21	0.40
1:C:300:LEU:HD13	1:C:307:ILE:HG12	2.02	0.40
1:D:164:VAL:CG2	4:D:469:HOH:O	2.69	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:7662:HOH:O	4:B:7691:HOH:O[2_555]	2.10	0.10
4:A:6430:HOH:O	4:B:7600:HOH:O[2_545]	2.18	0.02

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	391/392 (100%)	371 (95%)	20 (5%)	0	100	100
1	B	391/392 (100%)	373 (95%)	16 (4%)	2 (0%)	24	15
1	C	391/392 (100%)	372 (95%)	17 (4%)	2 (0%)	24	15
1	D	391/392 (100%)	375 (96%)	14 (4%)	2 (0%)	24	15
All	All	1564/1568 (100%)	1491 (95%)	67 (4%)	6 (0%)	30	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	350	ILE
1	C	3	PRO
1	B	291	GLY
1	D	3	PRO
1	C	350	ILE
1	B	350	ILE

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	279/278 (100%)	256 (92%)	23 (8%)	10	2
1	B	279/278 (100%)	258 (92%)	21 (8%)	12	3
1	C	279/278 (100%)	266 (95%)	13 (5%)	23	10
1	D	279/278 (100%)	261 (94%)	18 (6%)	15	4
All	All	1116/1112 (100%)	1041 (93%)	75 (7%)	15	4

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	LEU
1	A	89	CYS
1	A	110	ILE
1	A	149	THR
1	A	191	LYS
1	A	207	ARG
1	A	226	ASP
1	A	228	MET
1	A	230	LYS
1	A	237	LYS
1	A	258	LEU
1	A	263	GLU
1	A	272	LEU
1	A	276	VAL
1	A	298	LYS
1	A	322	GLN
1	A	328	LYS
1	A	332	TRP
1	A	339	VAL
1	A	358	LEU
1	A	359	ASN
1	A	367	ARG
1	A	392	LEU
1	B	39	LEU

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Mol	Chain	Res	Type
1	B	66	PRO
1	B	89	CYS
1	B	134	MET
1	B	191	LYS
1	B	207	ARG
1	B	220	ARG
1	B	228	MET
1	B	240	THR
1	B	257	LEU
1	B	263	GLU
1	B	272	LEU
1	B	276	VAL
1	B	298	LYS
1	B	322	GLN
1	B	328	LYS
1	B	332	TRP
1	B	339	VAL
1	B	353	SER
1	B	358	LEU
1	B	359	ASN
1	C	39	LEU
1	C	191	LYS
1	C	207	ARG
1	C	220	ARG
1	C	224	THR
1	C	260	SER
1	C	272	LEU
1	C	298	LYS
1	C	322	GLN
1	C	332	TRP
1	C	358	LEU
1	C	366	LYS
1	C	371	ARG
1	D	1	SER
1	D	39	LEU
1	D	89	CYS
1	D	134	MET
1	D	155	TYR
1	D	187	GLU
1	D	207	ARG
1	D	258	LEU
1	D	272	LEU

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Mol	Chain	Res	Type
1	D	276	VAL
1	D	298	LYS
1	D	322	GLN
1	D	339	VAL
1	D	353	SER
1	D	358	LEU
1	D	365	MET
1	D	367	ARG
1	D	371	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	78	GLN
1	A	175	GLN
1	A	184	ASN
1	B	78	GLN
1	B	175	GLN
1	B	184	ASN
1	B	221	HIS
1	C	78	GLN
1	C	101	GLN
1	C	175	GLN
1	C	184	ASN
1	C	221	HIS
1	C	316	ASN
1	D	78	GLN
1	D	102	GLN
1	D	184	ASN
1	D	316	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	5719	-	4,4,4	0.24	0	6,6,6	0.48	0
3	GOL	A	6394	-	5,5,5	0.48	0	5,5,5	1.30	0
3	GOL	B	7394	-	5,5,5	0.85	0	5,5,5	1.32	0
2	SO4	A	5720	-	4,4,4	0.20	0	6,6,6	0.35	0
2	SO4	B	5721	-	4,4,4	0.32	0	6,6,6	0.49	0
3	GOL	B	7393	-	5,5,5	0.70	0	5,5,5	1.28	0
2	SO4	A	5722	-	4,4,4	0.24	0	6,6,6	0.41	0
3	GOL	A	6393	-	5,5,5	0.58	0	5,5,5	0.89	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	7394	-	-	4/4/4/4	-
3	GOL	A	6394	-	-	4/4/4/4	-
3	GOL	A	6393	-	-	3/4/4/4	-
3	GOL	B	7393	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	6393	GOL	O1-C1-C2-C3
3	A	6394	GOL	O1-C1-C2-O2
3	A	6394	GOL	O1-C1-C2-C3
3	B	7394	GOL	O1-C1-C2-O2
3	B	7394	GOL	O1-C1-C2-C3
3	A	6394	GOL	C1-C2-C3-O3
3	B	7394	GOL	C1-C2-C3-O3
3	A	6393	GOL	O1-C1-C2-O2
3	B	7394	GOL	O2-C2-C3-O3
3	A	6394	GOL	O2-C2-C3-O3
3	A	6393	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	6394	GOL	3	0
3	B	7394	GOL	3	0
3	B	7393	GOL	4	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	392/392 (100%)	-1.26	0 100 100	3, 9, 26, 50	1 (0%)
1	B	392/392 (100%)	-1.27	0 100 100	3, 9, 26, 47	1 (0%)
1	C	392/392 (100%)	-1.04	0 100 100	2, 7, 24, 45	1 (0%)
1	D	392/392 (100%)	-0.97	0 100 100	2, 8, 28, 42	1 (0%)
All	All	1568/1568 (100%)	-1.13	0 100 100	2, 8, 26, 50	4 (0%)

There are no RSRZ outliers to report.

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
2	SO4	A	5722	5/5	0.95	0.09	74,75,77,77	0
3	GOL	A	6394	6/6	0.97	0.07	20,26,31,32	0
3	GOL	A	6393	6/6	0.98	0.05	19,26,28,31	0
2	SO4	B	5719	5/5	0.98	0.06	71,72,72,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	B	7394	6/6	0.98	0.07	14,24,30,33	0
2	SO4	B	5721	5/5	0.99	0.04	45,45,47,47	0
3	GOL	B	7393	6/6	0.99	0.05	13,20,22,23	0
2	SO4	A	5720	5/5	0.99	0.04	57,58,60,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.