Full wwPDB X-ray Structure Validation Report (i)

Nov 7, 2023 – 05:00 AM EST

PDB ID	:	1IGW
Title	:	Crystal Structure of the Isocitrate Lyase from the A219C mutant of Escherichia coli
Authors	:	Britton, K.L.; Abeysinghe, I.S.B.; Baker, P.J.; Barynin, V.; Diehl, P.; Langridge, S.J.; McFadden, B.A.; Sedelnikova, S.E.; Stillman, T.J.; Weeradechapon, K.; Rice, D.W.
Deposited on	:	2001-04-18
Resolution	:	2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

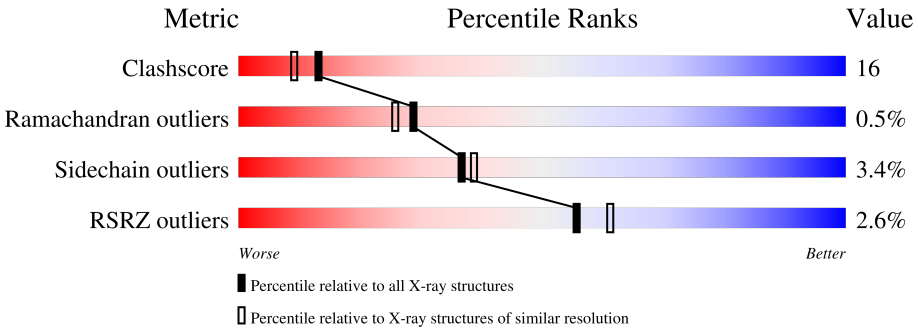
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	434	2% 73% 17% • 9%
1	B	434	3% 71% 21% • 5%
1	C	434	% 76% 18% • •
1	D	434	3% 69% 22% • 7%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HG	C	438	-	-	X	-
2	HG	D	439	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 13435 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 Isocitrate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	396	Total	C	N	O	S	0	0	0
			3066	1948	521	583	14			
1	B	411	Total	C	N	O	S	0	0	0
			3191	2028	542	608	13			
1	C	416	Total	C	N	O	S	0	0	0
			3223	2046	549	613	15			
1	D	405	Total	C	N	O	S	0	0	0
			3139	1995	531	600	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	219	CYS	ALA	engineered mutation	UNP P0A9G6
B	219	CYS	ALA	engineered mutation	UNP P0A9G6
C	219	CYS	ALA	engineered mutation	UNP P0A9G6
D	219	CYS	ALA	engineered mutation	UNP P0A9G6

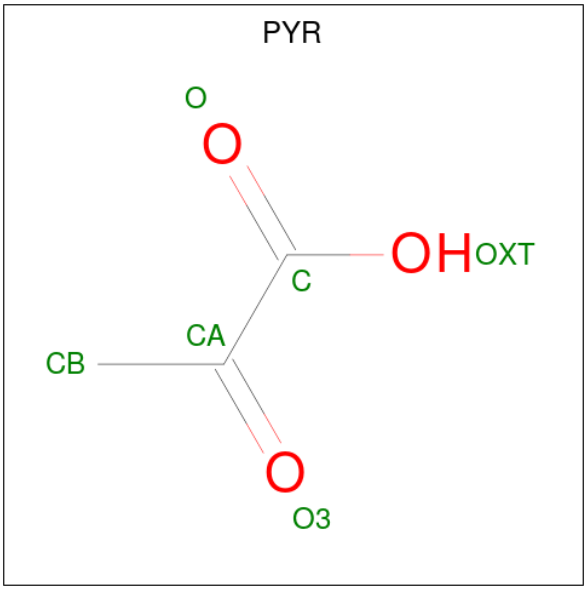
- Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Hg	0	0
			5	5		
2	B	5	Total	Hg	0	0
			5	5		
2	C	6	Total	Hg	0	0
			6	6		
2	D	5	Total	Hg	0	0
			5	5		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0

- Molecule 4 is PYRUVIC ACID (three-letter code: PYR) (formula: $C_3H_4O_3$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	188	Total O 188 188	0	0
5	B	190	Total O 190 190	0	0

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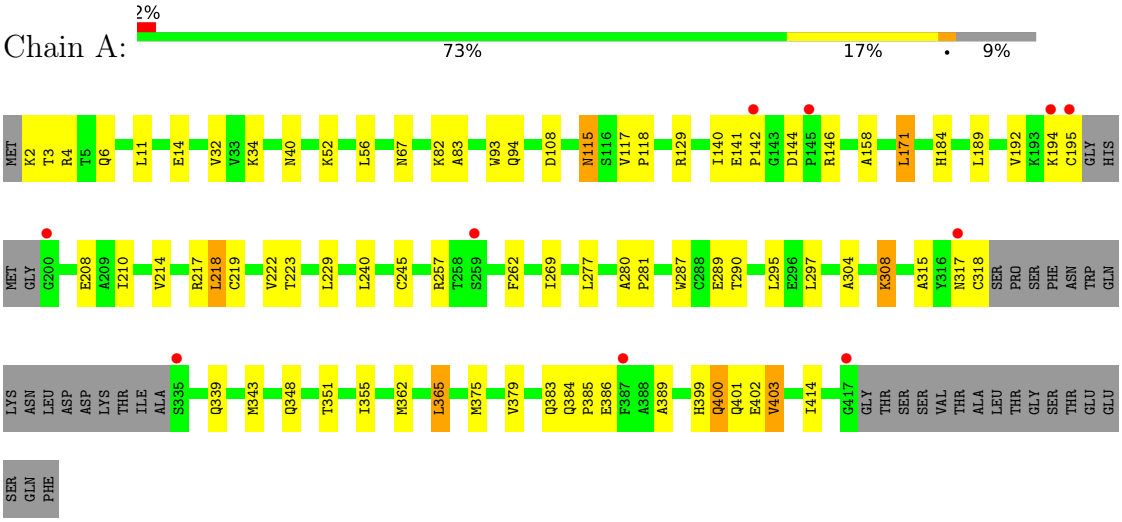
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	250	Total O 250 250	0	0
5	D	139	Total O 139 139	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

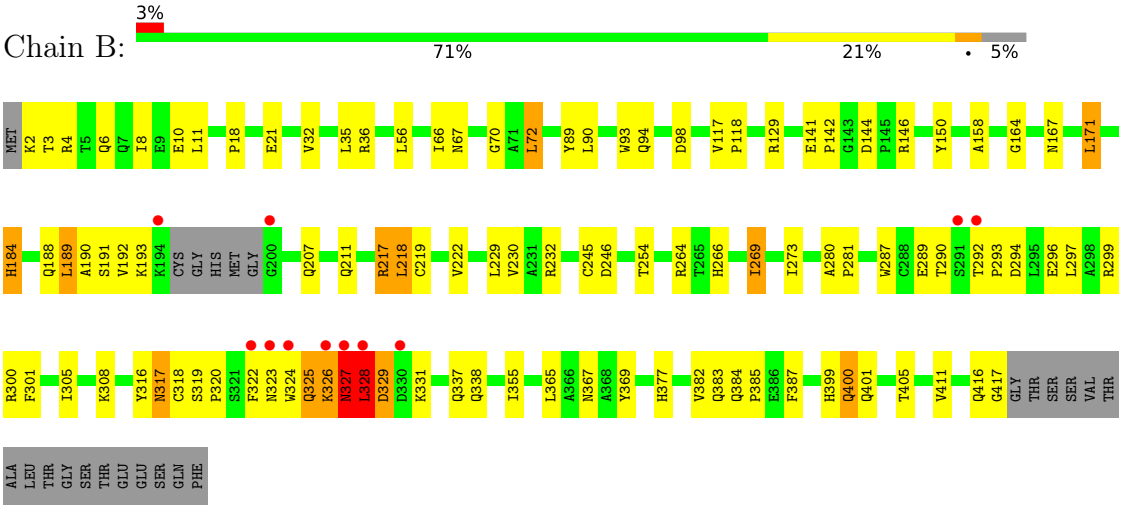
• Molecule 1: Isocitrate lyase

Chain A:



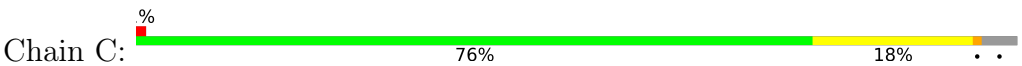
• Molecule 1: Isocitrate lyase

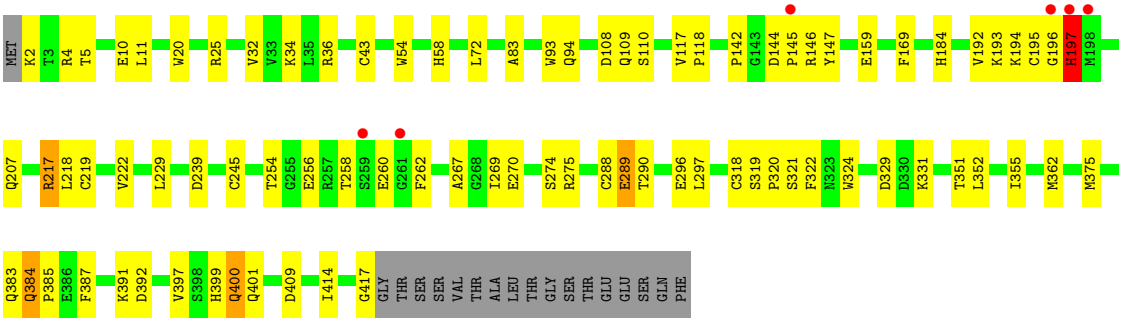
Chain B:



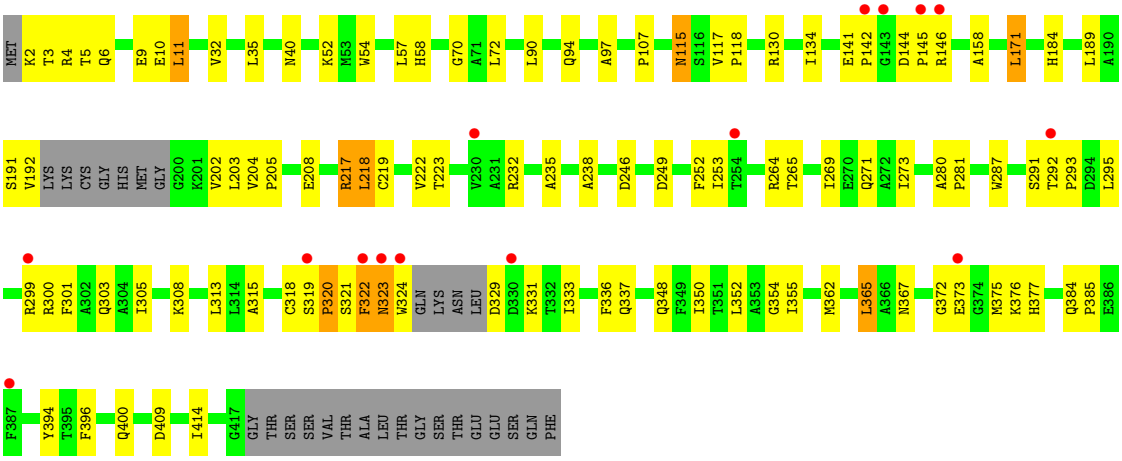
• Molecule 1: Isocitrate lyase

Chain C:





• Molecule 1: Isocitrate lyase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	88.65Å 88.65Å 199.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.10 33.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.10) 91.5 (33.24-2.10)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.84 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.235 0.171 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	22.7	Xtriage
Anisotropy	0.388	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 63.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -h,-k,l 0.035 for h,-h-k,-l 0.018 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	13435	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: HG, MG, PYR

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	0.40	0/3134	0.94	4/4251 (0.1%)
1	B	0.44	1/3264 (0.0%)	1.06	17/4430 (0.4%)
1	C	0.41	0/3298	0.98	7/4476 (0.2%)
1	D	0.36	0/3211	0.88	3/4360 (0.1%)
All	All	0.40	1/12907 (0.0%)	0.97	31/17517 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	405	THR	C-N	-7.21	1.20	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	217	ARG	NE-CZ-NH1	9.95	125.27	120.30
1	A	2	LYS	N-CA-C	-9.24	86.04	111.00
1	B	72	LEU	CA-CB-CG	8.78	135.49	115.30
1	D	217	ARG	NE-CZ-NH2	-8.73	115.94	120.30
1	B	217	ARG	NE-CZ-NH2	-7.34	116.63	120.30
1	B	327	ASN	N-CA-CB	7.32	123.77	110.60
1	A	184	HIS	CA-CB-CG	7.30	126.02	113.60
1	B	327	ASN	CA-C-O	7.16	135.14	120.10
1	C	184	HIS	CA-CB-CG	7.11	125.69	113.60
1	C	25	ARG	CD-NE-CZ	7.10	133.54	123.60
1	D	217	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	B	327	ASN	C-N-CA	6.85	138.83	121.70
1	B	369	TYR	CA-CB-CG	6.66	126.05	113.40
1	B	129	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	B	36	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	B	129	ARG	NE-CZ-NH2	-6.48	117.06	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	318	CYS	N-CA-CB	6.42	122.16	110.60
1	B	98	ASP	CB-CG-OD1	5.97	123.67	118.30
1	D	184	HIS	CA-CB-CG	5.91	123.65	113.60
1	B	184	HIS	CA-CB-CG	5.72	123.32	113.60
1	B	150	TYR	CB-CG-CD2	-5.70	117.58	121.00
1	B	217	ARG	CD-NE-CZ	5.61	131.46	123.60
1	C	275	ARG	NE-CZ-NH2	5.56	123.08	120.30
1	B	327	ASN	CA-C-N	-5.55	105.00	117.20
1	C	196	GLY	C-N-CA	5.47	135.38	121.70
1	A	129	ARG	CD-NE-CZ	5.46	131.25	123.60
1	A	129	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	B	399	HIS	CA-CB-CG	5.32	122.64	113.60
1	C	217	ARG	CD-NE-CZ	5.27	130.97	123.60
1	B	328	LEU	N-CA-C	-5.07	97.31	111.00
1	C	36	ARG	NE-CZ-NH2	-5.05	117.78	120.30

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	3066	0	2980	89	0
1	B	3191	0	3099	101	0
1	C	3223	0	3127	91	0
1	D	3139	0	3036	131	0
2	A	5	0	0	1	0
2	B	5	0	0	2	0
2	C	6	0	0	2	0
2	D	5	0	0	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	6	0	0	2	0
4	B	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	6	0	0	2	0
4	D	6	0	0	0	0
5	A	188	0	0	11	0
5	B	190	0	0	23	0
5	C	250	0	0	22	0
5	D	139	0	0	20	0
All	All	13435	0	12242	396	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 16.

All (396) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:293:PRO:HD3	1:D:322:PHE:CE1	1.53	1.43
1:D:293:PRO:HD3	1:D:322:PHE:CZ	1.56	1.39
1:D:409:ASP:HB3	5:D:1539:HOH:O	1.19	1.33
1:A:195:CYS:HB2	5:D:1565:HOH:O	1.30	1.25
1:B:416:GLN:OE1	1:C:195:CYS:SG	1.97	1.23
2:A:435:HG:HG	5:A:1445:HOH:O	1.15	1.20
1:A:290:THR:HG22	5:A:1558:HOH:O	1.35	1.20
1:B:325:GLN:O	1:B:327:ASN:N	1.76	1.18
1:D:204:VAL:HG13	1:D:205:PRO:HD2	1.34	1.10
1:B:296:GLU:OE1	1:B:299:ARG:NH1	1.85	1.08
1:C:289:GLU:HB3	5:C:1666:HOH:O	1.53	1.08
2:C:438:HG:HG	5:C:1640:HOH:O	1.29	1.08
1:D:292:THR:HA	1:D:322:PHE:HE1	1.10	1.06
1:D:292:THR:HA	1:D:322:PHE:CE1	1.92	1.04
1:D:293:PRO:CD	1:D:322:PHE:CE1	2.40	1.03
1:D:318:CYS:SG	1:D:320:PRO:HB3	2.01	1.01
1:B:323:ASN:O	1:B:326:LYS:HB2	1.60	1.00
2:C:438:HG:HG	5:C:1516:HOH:O	1.40	0.99
1:B:2:LYS:HE2	1:B:10:GLU:OE2	1.61	0.99
1:C:383:GLN:HG3	5:C:1621:HOH:O	1.63	0.98
1:D:293:PRO:CD	1:D:322:PHE:CZ	2.46	0.98
1:C:290:THR:HG22	5:C:1640:HOH:O	1.64	0.97
1:D:291:SER:O	1:D:322:PHE:CE1	2.18	0.96
1:A:245:CYS:SG	5:B:1609:HOH:O	2.24	0.96
1:C:387:PHE:CE2	5:C:1621:HOH:O	2.16	0.95
1:D:318:CYS:HB3	2:D:439:HG:HG	1.26	0.94
1:D:2:LYS:HE2	1:D:10:GLU:OE2	1.69	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:GLN:HE21	1:A:400:GLN:HA	1.34	0.92
1:B:329:ASP:OD2	1:B:331:LYS:CG	2.18	0.91
1:A:195:CYS:SG	5:A:1629:HOH:O	2.27	0.91
1:D:295:LEU:HD11	1:D:336:PHE:CE2	2.06	0.91
1:D:319:SER:HB2	1:D:321:SER:N	1.86	0.89
1:D:319:SER:N	5:D:1580:HOH:O	2.05	0.89
1:B:325:GLN:C	1:B:327:ASN:N	2.24	0.89
1:B:325:GLN:C	1:B:327:ASN:H	1.73	0.87
1:A:318:CYS:HA	5:A:1599:HOH:O	1.75	0.86
1:A:192:VAL:HG12	1:A:192:VAL:O	1.76	0.85
1:D:318:CYS:HG	1:D:324:TRP:HZ2	0.93	0.85
1:B:329:ASP:OD2	1:B:331:LYS:HG3	1.74	0.85
1:D:318:CYS:SG	5:D:1580:HOH:O	2.33	0.85
5:B:1617:HOH:O	1:C:375:MET:HE1	1.75	0.85
1:D:318:CYS:CB	2:D:439:HG:HG	1.86	0.84
1:D:322:PHE:CE2	5:D:1543:HOH:O	2.30	0.84
1:B:264:ARG:HD2	5:B:1567:HOH:O	1.76	0.84
1:B:325:GLN:NE2	1:C:383:GLN:OE1	2.10	0.83
1:D:52:LYS:HB2	1:D:52:LYS:NZ	1.92	0.83
1:B:323:ASN:O	1:B:326:LYS:CB	2.26	0.83
1:D:217:ARG:HD2	5:D:1491:HOH:O	1.78	0.82
1:B:318:CYS:SG	5:B:1617:HOH:O	2.36	0.82
1:C:2:LYS:HE2	1:C:10:GLU:OE2	1.79	0.81
1:D:204:VAL:CG1	1:D:205:PRO:HD2	2.10	0.81
1:C:296:GLU:HG3	5:C:1668:HOH:O	1.80	0.81
1:A:384:GLN:HB3	1:A:385:PRO:HD3	1.62	0.81
4:C:1446:PYR:CB	5:C:1635:HOH:O	2.29	0.81
1:D:318:CYS:SG	1:D:324:TRP:HZ2	2.04	0.80
1:B:401:GLN:HG3	5:B:1607:HOH:O	1.80	0.80
1:B:327:ASN:O	1:B:328:LEU:HG	1.80	0.79
1:A:144:ASP:OD2	1:A:146:ARG:CG	2.31	0.79
1:A:289:GLU:HG2	1:A:317:ASN:HD22	1.47	0.79
1:D:319:SER:HB2	1:D:320:PRO:CA	2.12	0.79
1:D:319:SER:HB2	1:D:321:SER:H	1.45	0.78
1:C:245:CYS:SG	5:C:1652:HOH:O	2.42	0.78
1:A:115:ASN:H	1:A:115:ASN:HD22	1.31	0.77
1:D:219:CYS:HG	2:D:436:HG:HG	1.27	0.77
1:D:94:GLN:NE2	1:D:355:ILE:H	1.82	0.77
1:D:3:THR:HB	1:D:6:GLN:HG3	1.65	0.77
1:C:207:GLN:HA	1:C:207:GLN:NE2	2.01	0.76
1:A:257:ARG:NH2	5:A:1607:HOH:O	2.20	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:THR:CG2	1:C:256:GLU:HG3	2.16	0.75
1:A:375:MET:HE2	1:D:324:TRP:CZ2	2.22	0.74
1:B:293:PRO:CD	5:B:1631:HOH:O	2.34	0.74
1:B:192:VAL:HG22	1:B:192:VAL:O	1.87	0.73
1:D:329:ASP:OD2	1:D:331:LYS:CG	2.36	0.73
1:A:400:GLN:HA	1:A:400:GLN:NE2	2.02	0.73
1:B:245:CYS:HG	2:B:437:HG:HG	0.72	0.73
1:D:158:ALA:HB1	1:D:171:LEU:HD13	1.70	0.73
1:C:254:THR:HG22	1:C:256:GLU:H	1.55	0.72
1:C:400:GLN:HA	1:C:400:GLN:HE21	1.55	0.72
1:A:144:ASP:OD2	1:A:146:ARG:HG2	1.89	0.72
1:A:94:GLN:NE2	1:A:355:ILE:H	1.87	0.72
1:D:319:SER:HB2	1:D:320:PRO:HA	1.71	0.72
1:A:192:VAL:HG11	1:A:262:PHE:CZ	2.26	0.71
1:C:289:GLU:CD	5:C:1666:HOH:O	2.27	0.71
1:B:323:ASN:HB3	1:B:326:LYS:HD2	1.72	0.71
1:B:384:GLN:HB3	1:B:385:PRO:HD3	1.71	0.71
1:C:351:THR:HG21	4:C:1446:PYR:CB	2.20	0.71
1:D:52:LYS:HB2	1:D:52:LYS:HZ2	1.53	0.71
1:D:299:ARG:O	1:D:303:GLN:HG3	1.90	0.71
1:B:217:ARG:HD2	5:B:1501:HOH:O	1.90	0.71
1:C:409:ASP:HB2	5:C:1683:HOH:O	1.90	0.70
1:D:308:LYS:HG3	5:D:1564:HOH:O	1.91	0.70
1:B:293:PRO:N	5:B:1631:HOH:O	2.24	0.70
1:B:2:LYS:CE	1:B:10:GLU:OE2	2.40	0.69
1:B:320:PRO:HB3	5:B:1617:HOH:O	1.91	0.69
1:D:204:VAL:HG13	1:D:205:PRO:CD	2.19	0.69
1:A:192:VAL:HG11	1:A:262:PHE:HZ	1.57	0.69
1:A:277:LEU:HD21	1:A:308:LYS:HE2	1.74	0.69
1:C:397:VAL:O	5:C:1569:HOH:O	2.11	0.69
1:B:320:PRO:HG3	5:B:1617:HOH:O	1.94	0.68
1:D:365:LEU:O	1:D:365:LEU:HD23	1.94	0.68
1:A:14:GLU:HG2	1:A:14:GLU:O	1.94	0.68
1:B:94:GLN:NE2	1:B:355:ILE:H	1.91	0.67
1:D:2:LYS:CE	1:D:10:GLU:OE2	2.42	0.67
1:A:144:ASP:OD2	1:A:146:ARG:HG3	1.92	0.67
5:A:1507:HOH:O	1:D:375:MET:SD	2.53	0.67
1:B:377:HIS:HE1	5:B:1595:HOH:O	1.77	0.67
1:D:269:ILE:O	1:D:273:ILE:HG13	1.95	0.67
1:D:295:LEU:CD1	1:D:336:PHE:CE2	2.78	0.67
1:D:329:ASP:OD2	1:D:331:LYS:HG3	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ASP:OD2	1:B:331:LYS:HG2	1.93	0.67
1:C:197:HIS:ND1	1:C:197:HIS:N	2.43	0.67
1:D:318:CYS:CA	5:D:1580:HOH:O	2.43	0.66
1:D:94:GLN:HE21	1:D:355:ILE:HG22	1.61	0.66
1:B:269:ILE:O	1:B:273:ILE:HG13	1.94	0.66
5:B:1617:HOH:O	1:C:375:MET:CE	2.40	0.66
1:D:329:ASP:OD2	1:D:331:LYS:HG2	1.95	0.65
1:A:365:LEU:HD23	1:A:365:LEU:O	1.96	0.65
1:D:291:SER:O	1:D:322:PHE:CZ	2.49	0.65
1:B:327:ASN:O	1:B:328:LEU:CG	2.43	0.65
1:C:329:ASP:OD2	1:C:331:LYS:CG	2.45	0.65
1:D:365:LEU:HD23	1:D:365:LEU:C	2.17	0.65
1:B:293:PRO:HD3	5:B:1631:HOH:O	1.94	0.64
1:A:195:CYS:CB	5:D:1565:HOH:O	2.08	0.64
1:D:373:GLU:O	1:D:376:LYS:HB3	1.97	0.64
1:D:319:SER:HB2	1:D:320:PRO:C	2.18	0.63
1:D:365:LEU:C	1:D:365:LEU:CD2	2.67	0.63
1:B:141:GLU:OE2	1:D:4:ARG:NE	2.28	0.63
1:D:204:VAL:CG1	1:D:208:GLU:HB2	2.29	0.63
5:B:1490:HOH:O	1:C:375:MET:SD	2.56	0.63
1:C:417:GLY:HA3	5:C:1618:HOH:O	1.98	0.63
1:A:192:VAL:O	1:A:192:VAL:CG1	2.46	0.62
1:B:319:SER:N	1:B:320:PRO:HA	2.13	0.62
1:D:158:ALA:HB1	1:D:171:LEU:CD1	2.28	0.62
1:C:289:GLU:CB	5:C:1666:HOH:O	2.27	0.62
1:B:325:GLN:O	1:B:327:ASN:CA	2.48	0.62
1:C:383:GLN:HG2	1:C:387:PHE:CE1	2.34	0.62
1:D:130:ARG:O	1:D:134:ILE:HG13	1.99	0.62
1:B:324:TRP:HZ3	1:B:328:LEU:HD12	1.63	0.62
1:C:192:VAL:HG12	1:C:192:VAL:O	1.99	0.62
1:C:322:PHE:HE2	1:C:324:TRP:CG	2.17	0.62
1:C:218:LEU:O	1:C:222:VAL:HG23	1.99	0.62
1:C:254:THR:HG22	1:C:256:GLU:HG3	1.82	0.62
1:D:115:ASN:HD22	1:D:115:ASN:H	1.47	0.62
1:A:277:LEU:HD11	1:A:308:LYS:HE3	1.82	0.61
1:C:319:SER:N	1:C:320:PRO:HA	2.15	0.61
1:A:287:TRP:HE1	1:A:317:ASN:CB	2.13	0.61
1:C:329:ASP:OD2	1:C:331:LYS:HG2	2.01	0.61
1:A:375:MET:HE2	1:D:324:TRP:CE2	2.36	0.60
1:C:322:PHE:CE2	1:C:324:TRP:CD1	2.89	0.60
1:B:325:GLN:O	1:B:326:LYS:C	2.38	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:218:LEU:O	1:D:222:VAL:HG23	2.02	0.60
1:A:277:LEU:CD2	1:A:308:LYS:HG2	2.32	0.59
1:B:67:ASN:HD22	1:B:337:GLN:HE22	1.50	0.59
1:B:327:ASN:O	1:B:328:LEU:CD2	2.51	0.59
1:A:117:VAL:N	1:A:118:PRO:CD	2.65	0.59
1:C:383:GLN:HG2	1:C:387:PHE:CZ	2.38	0.59
1:A:295:LEU:HD22	1:A:343:MET:CE	2.33	0.58
1:C:319:SER:OG	1:C:321:SER:N	2.37	0.58
1:A:3:THR:HB	1:A:6:GLN:NE2	2.19	0.58
1:D:191:SER:HB3	5:D:1577:HOH:O	2.04	0.58
1:A:195:CYS:SG	5:D:1565:HOH:O	2.56	0.58
1:A:365:LEU:C	1:A:365:LEU:CD2	2.71	0.57
1:A:141:GLU:HG2	1:C:4:ARG:HH21	1.70	0.57
1:A:217:ARG:HD2	5:A:1495:HOH:O	2.03	0.57
1:C:391:LYS:HG3	1:C:392:ASP:N	2.19	0.57
1:D:253:ILE:HD12	5:D:1547:HOH:O	2.05	0.57
1:D:322:PHE:CD2	5:D:1543:HOH:O	2.56	0.56
1:D:372:GLY:O	1:D:377:HIS:HE1	1.88	0.56
1:D:384:GLN:HB3	1:D:385:PRO:HD3	1.86	0.56
1:A:287:TRP:HB2	1:A:315:ALA:HB3	1.87	0.56
1:A:158:ALA:HB1	1:A:171:LEU:HD13	1.87	0.56
1:A:403:VAL:HG21	1:D:97:ALA:HB2	1.88	0.56
1:C:289:GLU:CG	5:C:1666:HOH:O	2.51	0.56
1:C:329:ASP:OD2	1:C:331:LYS:HG3	2.06	0.56
1:A:141:GLU:HB3	1:A:142:PRO:HD2	1.87	0.56
1:D:189:LEU:HB3	1:D:192:VAL:HG23	1.88	0.56
5:B:1453:HOH:O	1:C:375:MET:HE2	2.05	0.56
1:C:269:ILE:HD12	1:C:297:LEU:CD1	2.36	0.56
1:A:277:LEU:HD22	1:A:308:LYS:HG2	1.87	0.55
1:A:277:LEU:HD22	1:A:308:LYS:CG	2.37	0.55
1:B:117:VAL:N	1:B:118:PRO:CD	2.70	0.55
1:B:320:PRO:CB	5:B:1617:HOH:O	2.50	0.55
1:B:144:ASP:OD2	1:B:146:ARG:CG	2.54	0.55
1:C:147:TYR:O	1:C:147:TYR:CD2	2.59	0.55
1:D:141:GLU:HB3	1:D:142:PRO:HD2	1.88	0.55
1:A:218:LEU:O	1:A:222:VAL:HG23	2.06	0.55
1:B:219:CYS:HG	2:B:436:HG:HG	1.48	0.55
1:A:287:TRP:HE1	1:A:317:ASN:HB3	1.71	0.55
1:B:158:ALA:HB1	1:B:171:LEU:HD13	1.88	0.55
1:C:192:VAL:CG1	1:C:262:PHE:HZ	2.20	0.55
1:D:203:LEU:HG	1:D:238:ALA:CB	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:319:SER:CB	1:D:320:PRO:CA	2.81	0.54
1:A:11:LEU:HD11	1:A:32:VAL:HG11	1.90	0.54
1:C:207:GLN:HE21	1:C:207:GLN:CA	2.19	0.54
1:A:308:LYS:HD3	1:A:308:LYS:N	2.23	0.54
1:D:52:LYS:HB2	1:D:52:LYS:HZ3	1.72	0.54
1:D:52:LYS:NZ	1:D:52:LYS:CB	2.66	0.54
1:D:400:GLN:NE2	5:D:1569:HOH:O	2.40	0.54
1:B:141:GLU:HB3	1:B:142:PRO:HD2	1.88	0.54
1:D:362:MET:HE3	1:D:362:MET:HA	1.89	0.54
1:B:327:ASN:O	1:B:328:LEU:HD23	2.08	0.54
1:A:208:GLU:OE1	5:A:1606:HOH:O	2.18	0.53
1:A:295:LEU:HD22	1:A:343:MET:HE1	1.91	0.53
1:C:362:MET:HA	1:C:362:MET:HE3	1.91	0.53
1:D:204:VAL:CG1	1:D:205:PRO:CD	2.81	0.53
1:B:293:PRO:HA	1:B:316:TYR:OH	2.09	0.53
1:D:394:TYR:CZ	1:D:396:PHE:HB3	2.44	0.53
1:A:365:LEU:HD23	1:A:365:LEU:C	2.29	0.53
1:A:400:GLN:HE21	1:A:400:GLN:CA	2.06	0.53
1:C:94:GLN:NE2	1:C:355:ILE:H	2.07	0.52
1:D:219:CYS:O	1:D:223:THR:HG23	2.08	0.52
1:B:144:ASP:OD2	1:B:146:ARG:HG2	2.09	0.52
1:B:377:HIS:HD2	5:B:1547:HOH:O	1.90	0.52
1:A:414:ILE:HD11	1:C:34:LYS:CE	2.39	0.52
1:D:115:ASN:HD22	1:D:115:ASN:N	2.08	0.52
1:A:401:GLN:OE1	1:A:401:GLN:N	2.42	0.52
1:A:3:THR:H	1:A:6:GLN:NE2	2.08	0.52
1:A:3:THR:H	1:A:6:GLN:HE21	1.58	0.52
1:D:320:PRO:HA	5:D:1580:HOH:O	2.10	0.52
1:D:318:CYS:HA	5:D:1580:HOH:O	2.05	0.52
5:A:1507:HOH:O	1:D:375:MET:CG	2.57	0.52
1:C:269:ILE:HD12	1:C:297:LEU:HD12	1.92	0.52
1:A:52:LYS:O	1:A:56:LEU:HG	2.11	0.51
1:D:204:VAL:HG12	1:D:208:GLU:HB2	1.91	0.51
1:D:70:GLY:O	1:D:354:GLY:HA3	2.10	0.51
1:C:391:LYS:HE3	1:C:392:ASP:OD1	2.09	0.51
1:A:280:ALA:HB3	1:A:281:PRO:HD3	1.92	0.51
1:C:387:PHE:CZ	5:C:1621:HOH:O	2.50	0.51
1:D:232:ARG:HA	1:D:287:TRP:O	2.10	0.51
1:D:117:VAL:N	1:D:118:PRO:CD	2.74	0.51
1:B:232:ARG:CZ	1:B:289:GLU:HG3	2.40	0.50
1:D:144:ASP:O	1:D:146:ARG:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:400:GLN:HE21	1:D:400:GLN:HA	1.74	0.50
1:C:193:LYS:NZ	5:C:1670:HOH:O	2.36	0.50
1:B:280:ALA:HB3	1:B:281:PRO:HD3	1.93	0.50
1:C:319:SER:OG	1:C:320:PRO:C	2.50	0.50
1:D:189:LEU:HD13	1:D:192:VAL:HG23	1.93	0.50
1:B:190:ALA:HA	1:B:193:LYS:HG3	1.93	0.50
1:D:40:ASN:ND2	5:D:1489:HOH:O	2.45	0.50
1:B:323:ASN:CB	1:B:326:LYS:HD2	2.40	0.50
1:D:252:PHE:O	1:D:265:THR:HA	2.12	0.50
1:C:144:ASP:O	1:C:146:ARG:N	2.45	0.49
1:D:362:MET:O	1:D:362:MET:HE2	2.13	0.49
1:D:115:ASN:H	1:D:115:ASN:ND2	2.09	0.49
1:C:109:GLN:OE1	5:C:1558:HOH:O	2.17	0.49
1:B:383:GLN:HG2	1:B:387:PHE:CE1	2.48	0.49
1:C:43:CYS:HB3	5:C:1459:HOH:O	2.12	0.49
1:A:287:TRP:HE1	1:A:317:ASN:HB2	1.77	0.49
1:B:382:VAL:O	1:B:385:PRO:HD2	2.12	0.49
1:C:322:PHE:HE2	1:C:324:TRP:CD1	2.30	0.49
1:C:117:VAL:N	1:C:118:PRO:CD	2.75	0.49
1:B:287:TRP:HE1	1:B:317:ASN:ND2	2.10	0.48
1:A:82:LYS:HE3	1:A:402:GLU:OE2	2.13	0.48
1:B:325:GLN:OE1	1:B:326:LYS:N	2.46	0.48
1:C:383:GLN:CG	1:C:387:PHE:CZ	2.96	0.48
1:C:290:THR:HG22	5:C:1516:HOH:O	2.13	0.48
1:B:18:PRO:HA	1:B:21:GLU:HG3	1.96	0.48
1:B:56:LEU:HB3	1:B:66:ILE:HD11	1.95	0.48
1:B:292:THR:CA	5:B:1631:HOH:O	2.61	0.48
1:A:115:ASN:H	1:A:115:ASN:ND2	2.06	0.48
1:A:351:THR:HG21	4:A:1444:PYR:CB	2.43	0.48
1:D:3:THR:HG22	1:D:5:THR:N	2.28	0.48
1:D:322:PHE:O	1:D:323:ASN:HB2	2.13	0.48
1:A:362:MET:HE3	1:A:362:MET:HA	1.96	0.48
1:C:20:TRP:HZ3	1:C:217:ARG:NH1	2.11	0.48
1:C:270:GLU:HG3	5:C:1574:HOH:O	2.14	0.48
1:A:67:ASN:HA	1:A:348:GLN:O	2.14	0.47
1:A:399:HIS:CE1	1:A:400:GLN:HG2	2.50	0.47
1:B:292:THR:C	5:B:1631:HOH:O	2.49	0.47
1:A:219:CYS:O	1:A:223:THR:HG23	2.13	0.47
1:A:414:ILE:HD11	1:C:34:LYS:HE3	1.97	0.47
1:C:192:VAL:HG12	1:C:262:PHE:HZ	1.79	0.47
1:D:295:LEU:CD1	1:D:336:PHE:HE2	2.25	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:GLN:HA	1:C:207:GLN:HE21	1.72	0.47
1:A:277:LEU:HD11	1:A:308:LYS:CE	2.44	0.47
1:C:319:SER:H	1:C:320:PRO:HA	1.80	0.47
1:D:295:LEU:HD11	1:D:336:PHE:CD2	2.50	0.47
1:B:35:LEU:HD11	1:D:414:ILE:HD12	1.97	0.47
1:B:93:TRP:CZ2	1:C:399:HIS:HD2	2.32	0.47
1:B:320:PRO:HB2	1:B:324:TRP:CD1	2.49	0.47
1:B:417:GLY:O	5:B:1623:HOH:O	2.21	0.47
1:D:144:ASP:OD2	1:D:146:ARG:HG2	2.15	0.47
1:C:290:THR:CG2	5:C:1516:HOH:O	2.62	0.46
1:D:5:THR:O	1:D:9:GLU:HG3	2.15	0.46
1:A:269:ILE:HD12	1:A:297:LEU:HD12	1.98	0.46
1:C:194:LYS:HZ2	1:C:258:THR:CG2	2.28	0.46
1:B:411:VAL:HG22	1:D:35:LEU:HD13	1.96	0.46
1:D:372:GLY:O	1:D:373:GLU:HB2	2.15	0.46
1:A:194:LYS:HE2	1:A:240:LEU:HD11	1.97	0.46
1:D:144:ASP:OD2	1:D:146:ARG:CG	2.64	0.46
1:D:301:PHE:O	1:D:305:ILE:HG12	2.15	0.46
1:A:40:ASN:ND2	5:A:1487:HOH:O	2.48	0.46
1:C:319:SER:HA	1:C:351:THR:OG1	2.15	0.46
1:A:287:TRP:NE1	1:A:317:ASN:HB2	2.31	0.46
1:A:295:LEU:HD22	1:A:343:MET:HE3	1.97	0.46
1:B:383:GLN:HG2	1:B:387:PHE:HE1	1.81	0.45
1:C:239:ASP:HB3	1:C:267:ALA:HB2	1.98	0.45
1:C:400:GLN:HA	1:C:400:GLN:NE2	2.27	0.45
4:A:1444:PYR:CB	5:A:1570:HOH:O	2.64	0.45
1:C:288:CYS:O	1:C:290:THR:HG23	2.17	0.45
1:B:294:ASP:C	1:B:294:ASP:OD1	2.55	0.45
1:D:292:THR:CA	1:D:322:PHE:CE1	2.84	0.45
1:D:319:SER:CB	1:D:320:PRO:HA	2.38	0.45
1:B:3:THR:OG1	1:B:4:ARG:N	2.50	0.45
1:B:3:THR:CG2	1:B:6:GLN:HG3	2.47	0.45
1:B:190:ALA:CB	1:B:193:LYS:HE3	2.46	0.45
1:B:323:ASN:OD1	1:B:325:GLN:OE1	2.35	0.45
1:C:322:PHE:HE2	1:C:324:TRP:CD2	2.34	0.44
1:A:317:ASN:OD1	1:A:317:ASN:C	2.56	0.44
1:D:308:LYS:HA	5:D:1564:HOH:O	2.17	0.44
1:C:110:SER:OG	1:C:159:GLU:OE1	2.23	0.44
1:A:192:VAL:CG1	1:A:262:PHE:CZ	2.99	0.44
1:B:56:LEU:CB	1:B:66:ILE:HD11	2.48	0.44
1:B:93:TRP:HE1	1:C:400:GLN:HE22	1.64	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:ALA:HB1	1:D:367:ASN:HA	2.00	0.44
1:B:192:VAL:O	1:B:192:VAL:CG2	2.59	0.44
1:D:189:LEU:HD23	1:D:189:LEU:HA	1.75	0.44
1:D:235:ALA:HB1	1:D:271:GLN:HG2	2.00	0.44
1:C:94:GLN:NE2	1:C:352:LEU:HA	2.33	0.44
1:A:304:ALA:O	1:A:308:LYS:HE2	2.17	0.44
1:B:4:ARG:O	1:B:8:ILE:HG13	2.17	0.44
1:D:3:THR:HG22	1:D:5:THR:H	1.81	0.44
1:D:319:SER:N	1:D:320:PRO:HA	2.32	0.43
1:D:203:LEU:HG	1:D:238:ALA:HB1	2.01	0.43
1:D:246:ASP:HB3	1:D:249:ASP:OD2	2.18	0.43
1:D:280:ALA:HB3	1:D:281:PRO:HD3	2.00	0.43
1:A:141:GLU:HB3	1:A:142:PRO:CD	2.48	0.43
1:B:300:ARG:HG2	5:B:1615:HOH:O	2.19	0.43
1:B:400:GLN:HE22	1:C:93:TRP:HE1	1.66	0.43
1:D:57:LEU:HD13	1:D:313:LEU:HD13	2.00	0.43
1:D:158:ALA:CB	1:D:171:LEU:HD13	2.46	0.43
1:A:3:THR:CG2	1:A:4:ARG:N	2.81	0.43
1:A:289:GLU:CG	1:A:317:ASN:HD22	2.24	0.43
1:A:379:VAL:HA	1:A:383:GLN:HB2	2.00	0.43
1:A:403:VAL:HG13	1:D:107:PRO:HA	2.01	0.43
1:A:289:GLU:HG2	1:A:317:ASN:ND2	2.26	0.43
1:A:375:MET:SD	1:D:350:ILE:HD11	2.58	0.43
1:C:391:LYS:HE3	1:C:392:ASP:CG	2.39	0.43
1:D:320:PRO:O	1:D:320:PRO:CG	2.67	0.43
1:A:277:LEU:CD1	1:A:308:LYS:HE3	2.48	0.42
1:B:70:GLY:HA2	1:B:89:TYR:O	2.19	0.42
1:B:382:VAL:C	1:B:385:PRO:HD2	2.39	0.42
1:D:203:LEU:HD11	1:D:235:ALA:HA	2.00	0.42
1:B:329:ASP:OD2	1:B:331:LYS:CE	2.67	0.42
1:D:54:TRP:CZ2	1:D:58:HIS:CE1	3.07	0.42
1:D:192:VAL:HG11	1:D:202:VAL:HG21	2.01	0.42
1:D:287:TRP:HB2	1:D:315:ALA:HB3	2.02	0.42
1:B:190:ALA:C	1:B:192:VAL:H	2.21	0.42
1:C:93:TRP:CD1	1:C:108:ASP:HB2	2.54	0.42
1:B:188:GLN:OE1	1:B:193:LYS:HE2	2.20	0.42
1:B:207:GLN:O	1:B:211:GLN:HG3	2.20	0.42
1:C:11:LEU:HD11	1:C:32:VAL:HG11	2.02	0.42
1:D:11:LEU:HD11	1:D:32:VAL:HG11	2.01	0.42
1:B:292:THR:N	5:B:1631:HOH:O	2.53	0.42
1:C:144:ASP:C	1:C:146:ARG:H	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:384:GLN:N	1:C:385:PRO:CD	2.83	0.42
1:A:141:GLU:CG	1:C:4:ARG:HH21	2.31	0.42
1:B:189:LEU:HD22	1:B:191:SER:OG	2.20	0.42
1:C:194:LYS:NZ	1:C:258:THR:HB	2.35	0.42
1:B:144:ASP:OD2	1:B:146:ARG:HG3	2.20	0.41
1:B:189:LEU:O	1:B:192:VAL:HG12	2.20	0.41
1:D:269:ILE:HD13	1:D:300:ARG:HD2	2.01	0.41
1:D:308:LYS:CA	5:D:1564:HOH:O	2.68	0.41
1:C:54:TRP:CE2	1:C:58:HIS:CE1	3.08	0.41
1:B:164:GLY:H	1:B:167:ASN:HD22	1.68	0.41
1:B:301:PHE:O	1:B:305:ILE:HG12	2.20	0.41
1:D:191:SER:CB	5:D:1577:HOH:O	2.66	0.41
1:D:329:ASP:O	1:D:333:ILE:HG13	2.20	0.41
1:A:93:TRP:CD1	1:A:108:ASP:HB2	2.56	0.41
1:B:318:CYS:SG	1:B:320:PRO:HB3	2.60	0.41
1:B:319:SER:H	1:B:320:PRO:HA	1.82	0.41
1:D:337:GLN:HB3	1:D:348:GLN:OE1	2.20	0.41
1:A:210:ILE:O	1:A:214:VAL:HG23	2.21	0.41
1:B:11:LEU:HD11	1:B:32:VAL:HG11	2.03	0.41
1:B:246:ASP:OD1	5:B:1558:HOH:O	2.22	0.41
1:D:384:GLN:N	1:D:385:PRO:CD	2.83	0.41
1:C:169:PHE:CE1	1:C:219:CYS:HB2	2.55	0.41
1:A:277:LEU:HD22	1:A:308:LYS:HG3	2.03	0.41
1:D:384:GLN:CB	1:D:385:PRO:HD3	2.51	0.41
1:A:386:GLU:O	1:A:389:ALA:HB3	2.20	0.41
1:B:141:GLU:HB3	1:B:142:PRO:CD	2.51	0.41
1:B:218:LEU:O	1:B:222:VAL:HG23	2.21	0.41
1:D:2:LYS:NZ	1:D:10:GLU:OE2	2.53	0.41
1:D:90:LEU:HD22	1:D:117:VAL:HG13	2.02	0.41
1:D:94:GLN:NE2	1:D:352:LEU:HA	2.36	0.41
1:B:290:THR:H	1:B:317:ASN:HD21	1.69	0.40
1:B:184:HIS:HB3	1:B:230:VAL:HB	2.02	0.40
1:C:142:PRO:HA	1:C:147:TYR:CD1	2.57	0.40
1:C:192:VAL:HG11	1:C:262:PHE:HZ	1.86	0.40
1:C:194:LYS:HE3	1:C:260:GLU:OE2	2.21	0.40
1:A:414:ILE:HD11	1:C:34:LYS:HE2	2.02	0.40
1:B:324:TRP:HZ3	1:B:328:LEU:CD1	2.30	0.40
1:A:34:LYS:HE2	1:C:414:ILE:HD11	2.03	0.40
1:B:254:THR:HG22	1:B:266:HIS:CE1	2.56	0.40
1:B:319:SER:N	1:B:320:PRO:CA	2.83	0.40
1:B:367:ASN:HA	1:C:83:ALA:HB1	2.04	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	390/434 (90%)	375 (96%)	14 (4%)	1 (0%)	41	41
1	B	407/434 (94%)	390 (96%)	15 (4%)	2 (0%)	29	26
1	C	414/434 (95%)	401 (97%)	11 (3%)	2 (0%)	29	26
1	D	399/434 (92%)	381 (96%)	15 (4%)	3 (1%)	19	15
All	All	1610/1736 (93%)	1547 (96%)	55 (3%)	8 (0%)	29	26

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	326	LYS
1	C	197	HIS
1	D	322	PHE
1	D	323	ASN
1	B	328	LEU
1	C	145	PRO
1	D	145	PRO
1	A	140	ILE

5.3.2 Protein sidechains (i)

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	312/344 (91%)	302 (97%)	10 (3%)	39	41
1	B	326/344 (95%)	309 (95%)	17 (5%)	23	21
1	C	329/344 (96%)	320 (97%)	9 (3%)	44	48
1	D	320/344 (93%)	312 (98%)	8 (2%)	47	52
All	All	1287/1376 (94%)	1243 (97%)	44 (3%)	37	39

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	ASN
1	A	171	LEU
1	A	189	LEU
1	A	218	LEU
1	A	229	LEU
1	A	308	LYS
1	A	339	GLN
1	A	365	LEU
1	A	400	GLN
1	A	403	VAL
1	B	72	LEU
1	B	90	LEU
1	B	171	LEU
1	B	189	LEU
1	B	218	LEU
1	B	229	LEU
1	B	269	ILE
1	B	297	LEU
1	B	308	LYS
1	B	317	ASN
1	B	322	PHE
1	B	325	GLN
1	B	327	ASN
1	B	329	ASP
1	B	338	GLN
1	B	365	LEU
1	B	400	GLN
1	C	5	THR
1	C	72	LEU
1	C	197	HIS
1	C	229	LEU
1	C	274	SER
1	C	289	GLU

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Mol	Chain	Res	Type
1	C	384	GLN
1	C	400	GLN
1	C	401	GLN
1	D	11	LEU
1	D	72	LEU
1	D	115	ASN
1	D	171	LEU
1	D	218	LEU
1	D	264	ARG
1	D	320	PRO
1	D	365	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	40	ASN
1	A	94	GLN
1	A	115	ASN
1	A	266	HIS
1	A	337	GLN
1	A	339	GLN
1	A	400	GLN
1	B	40	ASN
1	B	94	GLN
1	B	167	ASN
1	B	266	HIS
1	B	303	GLN
1	B	317	ASN
1	B	337	GLN
1	B	377	HIS
1	B	399	HIS
1	B	400	GLN
1	B	416	GLN
1	C	40	ASN
1	C	94	GLN
1	C	109	GLN
1	C	207	GLN
1	C	303	GLN
1	C	399	HIS
1	C	400	GLN
1	D	40	ASN

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Mol	Chain	Res	Type
1	D	94	GLN
1	D	115	ASN
1	D	377	HIS
1	D	399	HIS
1	D	400	GLN

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

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 29 ligands modelled in this entry, 25 are monoatomic - leaving 4 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	$\#\|Z\| > 2$
4	PYR	A	1444	3	5,5,5	3.93	1 (20%)	3,6,6	5.78	1 (33%)
4	PYR	D	1447	3	5,5,5	3.82	1 (20%)	3,6,6	2.90	1 (33%)
4	PYR	C	1446	3	5,5,5	4.25	1 (20%)	3,6,6	1.58	1 (33%)
4	PYR	B	1445	3	5,5,5	0.97	0	3,6,6	2.98	2 (66%)

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
4	PYR	A	1444	3	-	2/4/4/4	-
4	PYR	D	1447	3	-	2/4/4/4	-
4	PYR	C	1446	3	-	0/4/4/4	-
4	PYR	B	1445	3	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1446	PYR	O3-CA	9.25	1.43	1.23
4	A	1444	PYR	O3-CA	8.45	1.41	1.23
4	D	1447	PYR	O3-CA	8.21	1.41	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	1444	PYR	O3-CA-CB	9.96	141.82	119.73
4	D	1447	PYR	O3-CA-CB	-4.76	109.19	119.73
4	B	1445	PYR	O3-CA-CB	-4.66	109.41	119.73
4	C	1446	PYR	OXT-C-CA	2.67	121.28	113.97
4	B	1445	PYR	OXT-C-CA	2.18	119.93	113.97

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1444	PYR	O-C-CA-CB
4	A	1444	PYR	OXT-C-CA-CB
4	D	1447	PYR	O-C-CA-CB
4	D	1447	PYR	OXT-C-CA-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1444	PYR	2	0
4	C	1446	PYR	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	405:THR	C	406:GLY	N	1.20

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, 95^{th} 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	396/434 (91%)	-0.34	10 (2%)	57	62	14, 27, 56, 86	0
1	B	411/434 (94%)	-0.43	11 (2%)	54	60	12, 22, 56, 83	0
1	C	416/434 (95%)	-0.47	6 (1%)	75	78	11, 22, 49, 85	0
1	D	405/434 (93%)	0.01	15 (3%)	41	48	16, 34, 73, 94	0
All	All	1628/1736 (93%)	-0.31	42 (2%)	56	61	11, 26, 62, 94	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	322	PHE	8.0
1	C	197	HIS	5.7
1	B	291	SER	4.4
1	B	326	LYS	4.1
1	C	145	PRO	4.0
1	D	145	PRO	3.9
1	B	322	PHE	3.6
1	D	319	SER	3.6
1	B	328	LEU	3.5
1	D	142	PRO	3.4
1	D	324	TRP	3.3
1	D	146	ARG	3.1
1	B	330	ASP	2.9
1	D	387	PHE	2.8
1	B	327	ASN	2.8
1	D	323	ASN	2.8
1	A	417	GLY	2.8
1	B	324	TRP	2.7
1	A	195	CYS	2.7
1	A	387	PHE	2.6
1	A	259	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	323	ASN	2.5
1	D	373	GLU	2.5
1	D	254	THR	2.5
1	C	259	SER	2.5
1	D	330	ASP	2.4
1	B	194	LYS	2.4
1	B	292	THR	2.4
1	A	335	SER	2.3
1	C	261	GLY	2.2
1	B	200	GLY	2.2
1	C	196	GLY	2.2
1	D	143	GLY	2.2
1	A	145	PRO	2.1
1	D	299	ARG	2.1
1	D	230	VAL	2.1
1	A	142	PRO	2.1
1	A	317	ASN	2.1
1	A	194	LYS	2.1
1	C	198	MET	2.0
1	D	292	THR	2.0
1	A	200	GLY	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 monosaccharides 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, 95^{th} 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
3	MG	D	441	1/1	0.88	0.12	27,27,27,27	0
3	MG	A	441	1/1	0.91	0.08	26,26,26,26	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	PYR	B	1445	6/6	0.96	0.09	7,14,17,17	0
4	PYR	C	1446	6/6	0.96	0.07	12,16,18,18	0
2	HG	D	437	1/1	0.97	0.04	43,43,43,43	1
4	PYR	A	1444	6/6	0.97	0.07	19,25,30,34	0
2	HG	D	438	1/1	0.97	0.03	55,55,55,55	0
2	HG	A	437	1/1	0.97	0.06	33,33,33,33	1
2	HG	C	437	1/1	0.98	0.05	43,43,43,43	1
2	HG	D	439	1/1	0.98	0.05	45,45,45,45	1
2	HG	B	437	1/1	0.98	0.06	39,39,39,39	1
3	MG	C	441	1/1	0.98	0.03	21,21,21,21	0
4	PYR	D	1447	6/6	0.98	0.12	25,26,28,28	0
2	HG	C	436	1/1	0.99	0.02	39,39,39,39	0
2	HG	B	436	1/1	0.99	0.03	37,37,37,37	0
3	MG	B	441	1/1	0.99	0.10	18,18,18,18	0
2	HG	C	438	1/1	0.99	0.01	30,30,30,30	0
2	HG	C	440	1/1	0.99	0.03	45,45,45,45	1
2	HG	D	435	1/1	0.99	0.05	31,31,31,31	0
2	HG	D	436	1/1	0.99	0.02	32,32,32,32	0
2	HG	A	438	1/1	0.99	0.03	37,37,37,37	0
2	HG	C	435	1/1	0.99	0.05	30,30,30,30	0
2	HG	B	438	1/1	1.00	0.04	37,37,37,37	0
2	HG	B	439	1/1	1.00	0.04	37,37,37,37	0
2	HG	A	439	1/1	1.00	0.03	40,40,40,40	1
2	HG	B	435	1/1	1.00	0.04	31,31,31,31	0
2	HG	A	435	1/1	1.00	0.03	31,31,31,31	0
2	HG	A	436	1/1	1.00	0.04	38,38,38,38	0
2	HG	C	439	1/1	1.00	0.01	24,24,24,24	0

6.5 Other polymers (i)

There are no such residues in this entry.