



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:33 AM UTC

PDB ID : 1W8Q / pdb_00001w8q
Title : Crystal Structure of the DD-Transpeptidase-carboxypeptidase from Actinomadura R39
Authors : Sauvage, E.; Herman, R.; Petrella, S.; Duez, C.; Frere, J.M.; Charlier, P.
Deposited on : 2004-09-24
Resolution : 2.85 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

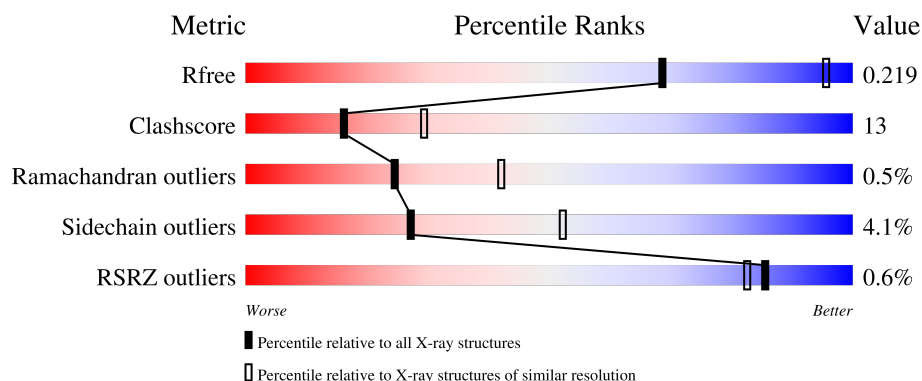
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	489	 72% 21% . .
1	B	489	 64% 30% . 5%
1	C	489	 67% 27% . 5%
1	D	489	 69% 25% . .

2 Entry composition [i](#)

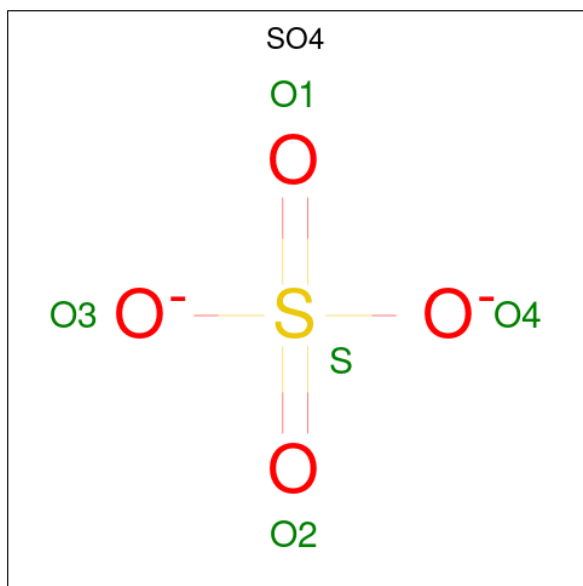
There are 4 unique types of molecules in this entry. The entry contains 13592 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-ALANYL-D-ALANINE CARBOXYPEPTIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	1
			3353	2076	565	706	6			
1	B	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	C	466	Total	C	N	O	S	0	0	1
			3344	2071	564	703	6			
1	D	467	Total	C	N	O	S	0	0	1
			3353	2076	565	706	6			

- 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		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0

- Molecule 3 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	5	Total Co 5 5	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total 4	Co 4	0	0
3	C	4	Total 4	Co 4	0	0
3	D	5	Total 5	Co 5	0	0

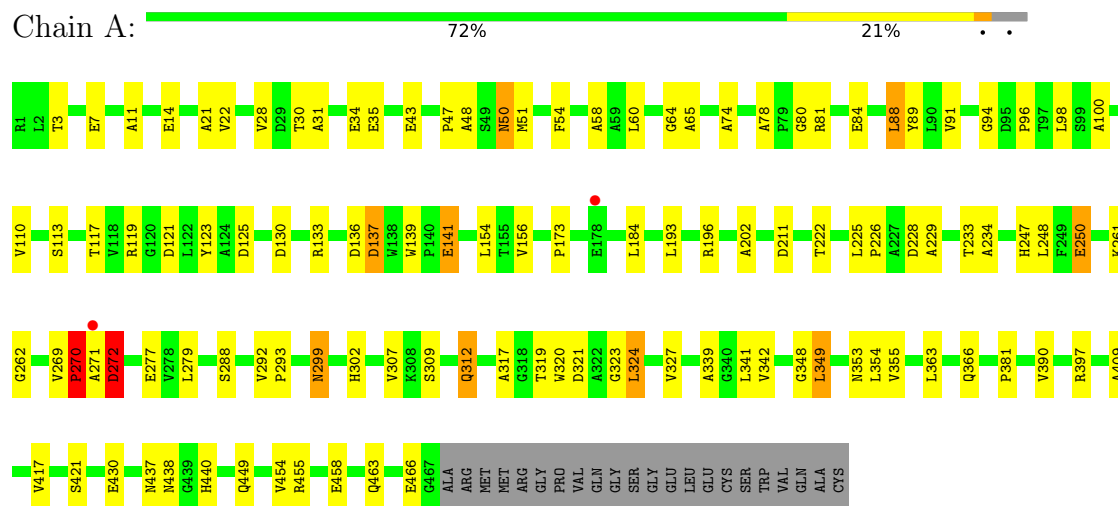
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total 25	O 25	0	0
4	B	18	Total 18	O 18	0	0
4	C	14	Total 14	O 14	0	0
4	D	23	Total 23	O 23	0	0

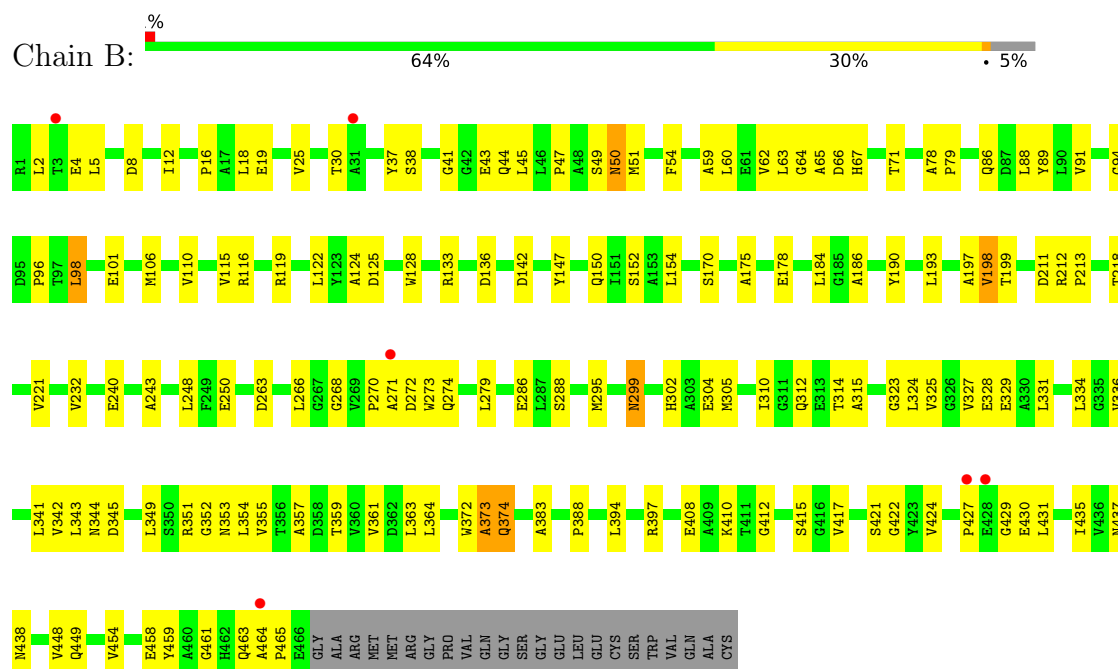
3 Residue-property plots

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: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

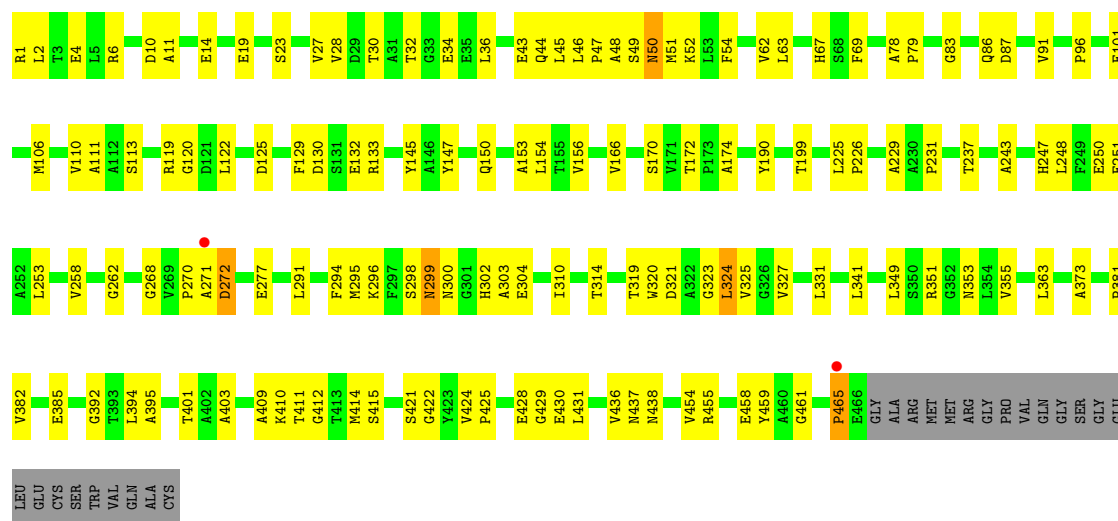


• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE



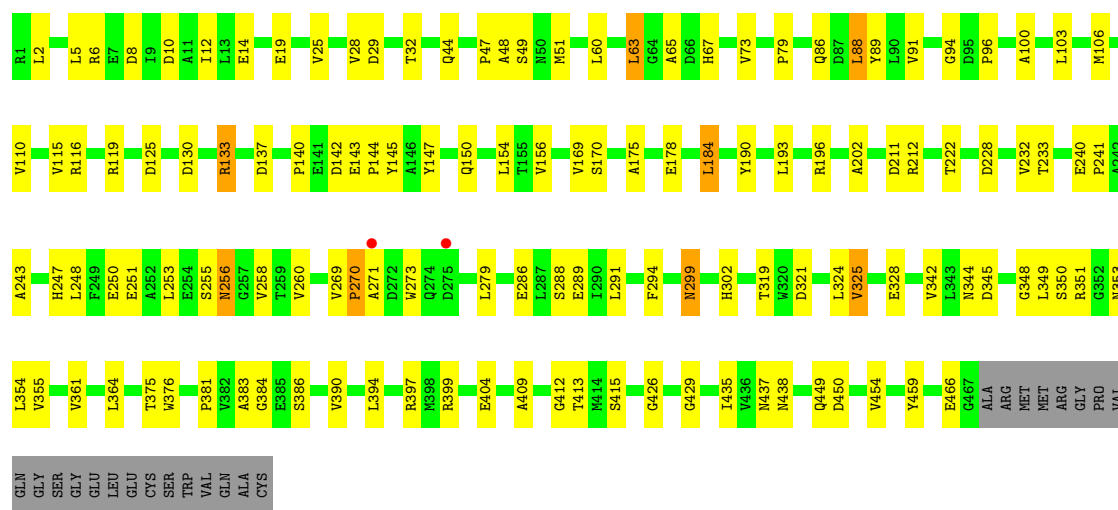
• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

Chain C:  67% 27% 5%



• Molecule 1: D-ALANYL-D-ALANINE CARBOXYPEPTIDASE

Chain D:  69% 25% 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	104.56Å 94.37Å 106.96Å 90.00° 94.05° 90.00°	Depositor
Resolution (Å)	19.94 – 2.85 19.94 – 2.85	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.94-2.85) 98.1 (19.94-2.85)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.86Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.197 , 0.244 (Not available) , 0.219	Depositor DCC
R_{free} test set	2391 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	46.2	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.027 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13592	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.99% 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, CO

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	0.43	1/3412 (0.0%)	0.93	9/4668 (0.2%)
1	B	0.41	1/3403 (0.0%)	0.92	3/4656 (0.1%)
1	C	0.40	1/3403 (0.0%)	0.91	3/4656 (0.1%)
1	D	0.45	1/3412 (0.0%)	0.93	7/4668 (0.1%)
All	All	0.42	4/13630 (0.0%)	0.92	22/18648 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	466	GLU	C-N	-5.78	1.25	1.33
1	B	465	PRO	C-N	-5.50	1.25	1.33
1	C	465	PRO	C-N	-5.45	1.25	1.33
1	A	466	GLU	C-N	-5.35	1.25	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	269	VAL	CA-C-N	7.79	128.51	119.47
1	D	269	VAL	C-N-CA	7.79	128.51	119.47
1	A	269	VAL	CA-C-N	6.47	127.93	119.84
1	A	269	VAL	C-N-CA	6.47	127.93	119.84
1	C	272	ASP	N-CA-C	-6.39	105.47	113.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	314	THR	N-CA-C	6.19	119.22	111.24
1	B	136	ASP	N-CA-C	5.98	118.28	111.11
1	A	390	VAL	N-CA-C	-5.63	107.39	112.12
1	D	133	ARG	N-CA-C	5.56	118.27	111.82
1	B	263	ASP	N-CA-C	5.47	117.59	110.53
1	A	272	ASP	N-CA-C	-5.43	105.52	111.82
1	A	100	ALA	N-CA-C	-5.41	105.55	111.82
1	A	91	VAL	N-CA-C	5.40	115.36	107.37
1	A	133	ARG	N-CA-C	5.29	117.13	111.36
1	D	116	ARG	N-CA-C	-5.25	108.14	114.75
1	D	325	VAL	N-CA-C	-5.18	105.49	110.72
1	D	390	VAL	N-CA-C	-5.18	107.10	111.56
1	D	270	PRO	N-CA-C	5.17	121.30	113.81
1	B	62	VAL	N-CA-C	5.12	116.46	110.62
1	A	137	ASP	N-CA-C	5.09	118.32	112.57
1	C	395	ALA	N-CA-C	5.05	116.59	111.14
1	A	309	SER	N-CA-C	-5.00	105.91	111.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	145	TYR	Sidechain

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	3353	0	3201	68	0
1	B	3344	0	3195	102	0
1	C	3344	0	3195	95	0
1	D	3353	0	3201	82	0
2	A	25	0	0	0	0
2	B	25	0	0	0	0
2	C	25	0	0	0	0
2	D	25	0	0	0	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	5	0	0	0	0
4	A	25	0	0	0	0
4	B	18	0	0	1	0
4	C	14	0	0	0	0
4	D	23	0	0	0	0
All	All	13592	0	12792	345	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 13.

All (345) 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:226:PRO:HG2	1:A:229:ALA:HB2	1.33	1.06
1:C:226:PRO:HG2	1:C:229:ALA:HB2	1.41	1.02
1:D:47:PRO:HG3	1:D:51:MET:HE2	1.46	0.97
1:B:65:ALA:HB1	1:B:288:SER:HB3	1.57	0.84
1:C:225:LEU:HD12	1:C:226:PRO:HD2	1.59	0.84
1:C:47:PRO:HG3	1:C:51:MET:HE2	1.62	0.82
1:A:47:PRO:HG3	1:A:51:MET:HE2	1.61	0.80
1:D:150:GLN:HE22	1:D:240:GLU:H	1.29	0.80
1:A:51:MET:HE3	1:A:353:ASN:HB3	1.65	0.78
1:C:130:ASP:HB2	1:C:319:THR:HG22	1.66	0.77
1:B:299:ASN:HD22	1:B:299:ASN:C	1.95	0.74
1:B:424:VAL:HB	1:B:431:LEU:HB2	1.71	0.73
1:C:296:LYS:HA	1:C:381:PRO:HG3	1.71	0.72
1:A:270:PRO:C	1:A:272:ASP:H	1.97	0.71
1:A:299:ASN:C	1:A:299:ASN:HD22	1.96	0.71
1:B:125:ASP:OD2	1:B:268:GLY:HA2	1.91	0.71
1:B:150:GLN:NE2	1:B:240:GLU:H	1.90	0.70
1:B:351:ARG:NH2	1:B:415:SER:HB3	2.07	0.69
1:C:270:PRO:C	1:C:272:ASP:H	1.98	0.69
1:A:342:VAL:HB	1:A:354:LEU:HB2	1.75	0.69
1:B:51:MET:HE3	1:B:353:ASN:HB3	1.75	0.68
1:D:397:ARG:HH12	1:D:449:GLN:HE21	1.41	0.68
1:A:51:MET:CE	1:A:353:ASN:HB3	2.24	0.68
1:D:299:ASN:HD22	1:D:302:HIS:H	1.40	0.67
1:D:256:ASN:N	1:D:256:ASN:HD22	1.93	0.67
1:B:397:ARG:HH12	1:B:449:GLN:HE21	1.41	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:LEU:HD21	1:C:36:LEU:HD22	1.76	0.67
1:D:140:PRO:HA	1:D:143:GLU:HG3	1.76	0.67
1:B:47:PRO:HG3	1:B:51:MET:HE2	1.78	0.66
1:C:129:PHE:HB2	1:C:153:ALA:HB2	1.77	0.66
1:B:12:ILE:HG22	1:B:448:VAL:HG13	1.78	0.65
1:D:345:ASP:OD1	1:D:350:SER:HB3	1.96	0.65
1:D:437:ASN:C	1:D:438:ASN:HD22	2.05	0.65
1:A:397:ARG:HH12	1:A:449:GLN:HE21	1.44	0.65
1:C:125:ASP:CG	1:C:268:GLY:HA2	2.22	0.65
1:D:196:ARG:HB2	1:D:222:THR:HG22	1.80	0.64
1:D:450:ASP:O	1:D:454:VAL:HG23	1.98	0.64
1:B:344:ASN:HD22	1:B:352:GLY:HA3	1.62	0.64
1:C:323:GLY:O	1:C:327:VAL:HG23	1.97	0.63
1:B:116:ARG:NH2	1:B:116:ARG:HB3	2.13	0.63
1:B:150:GLN:HE22	1:B:240:GLU:H	1.47	0.63
1:B:175:ALA:HB3	1:B:178:GLU:HG3	1.79	0.63
1:D:270:PRO:O	1:D:271:ALA:HB3	1.96	0.63
1:D:342:VAL:HB	1:D:354:LEU:HB2	1.79	0.63
1:C:2:LEU:HB3	1:C:6:ARG:HH22	1.64	0.63
1:D:299:ASN:ND2	1:D:302:HIS:H	1.97	0.62
1:C:2:LEU:HB3	1:C:6:ARG:NH2	2.14	0.62
1:A:437:ASN:C	1:A:438:ASN:HD22	2.07	0.62
1:B:106:MET:O	1:B:110:VAL:HG23	1.99	0.62
1:D:65:ALA:HB1	1:D:288:SER:HB3	1.81	0.62
1:D:130:ASP:HB3	1:D:319:THR:HG22	1.81	0.62
1:D:51:MET:HE3	1:D:353:ASN:HB3	1.81	0.61
1:D:94:GLY:O	1:D:96:PRO:HD3	2.00	0.61
1:C:62:VAL:HG11	1:C:310:ILE:HG23	1.82	0.61
1:D:156:VAL:HG21	1:D:248:LEU:HD12	1.83	0.61
1:B:343:LEU:N	1:B:343:LEU:HD12	2.16	0.61
1:D:348:GLY:HA2	1:D:353:ASN:ND2	2.14	0.61
1:B:323:GLY:O	1:B:327:VAL:HG23	2.01	0.61
1:B:454:VAL:O	1:B:458:GLU:HG3	2.01	0.59
1:C:437:ASN:C	1:C:438:ASN:HD22	2.10	0.59
1:D:384:GLY:HA3	1:D:404:GLU:HG3	1.82	0.59
1:D:255:SER:C	1:D:256:ASN:HD22	2.11	0.59
1:C:174:ALA:O	1:C:199:THR:HG21	2.03	0.59
1:D:150:GLN:NE2	1:D:240:GLU:H	1.99	0.58
1:C:87:ASP:OD1	1:C:120:GLY:HA3	2.04	0.58
1:D:383:ALA:HA	1:D:394:LEU:HB3	1.84	0.58
1:A:397:ARG:HH12	1:A:449:GLN:NE2	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:GLN:HA	1:B:355:VAL:O	2.04	0.58
1:A:226:PRO:HG2	1:A:229:ALA:CB	2.22	0.58
1:D:386:SER:HB3	1:D:399:ARG:NH2	2.19	0.58
1:A:319:THR:HG22	1:A:321:ASP:H	1.68	0.57
1:C:424:VAL:HB	1:C:431:LEU:HB2	1.86	0.57
1:C:106:MET:O	1:C:110:VAL:HG23	2.03	0.57
1:A:454:VAL:O	1:A:458:GLU:HG3	2.04	0.57
1:D:63:LEU:O	1:D:67:HIS:HB2	2.05	0.57
1:D:351:ARG:HH12	1:D:415:SER:HB3	1.70	0.57
1:D:47:PRO:CG	1:D:51:MET:HE2	2.27	0.57
1:A:30:THR:HB	1:A:430:GLU:CG	2.34	0.56
1:C:10:ASP:O	1:C:14:GLU:HG2	2.05	0.56
1:A:156:VAL:HG21	1:A:248:LEU:HD12	1.85	0.56
1:A:21:ALA:HB2	1:A:440:HIS:N	2.20	0.56
1:C:27:VAL:HG12	1:C:36:LEU:HD12	1.88	0.56
1:B:51:MET:CE	1:B:353:ASN:HB3	2.36	0.56
1:B:124:ALA:HB3	1:B:266:LEU:HD23	1.86	0.56
1:D:345:ASP:OD1	1:D:353:ASN:ND2	2.38	0.56
1:A:270:PRO:O	1:A:271:ALA:HB3	2.06	0.56
1:A:320:TRP:O	1:A:324:LEU:HB2	2.06	0.55
1:B:91:VAL:HG22	1:B:125:ASP:HB3	1.88	0.55
1:B:170:SER:HA	1:B:232:VAL:O	2.06	0.55
1:C:79:PRO:HB2	1:C:83:GLY:HA2	1.88	0.55
1:B:271:ALA:C	1:B:273:TRP:H	2.14	0.55
1:C:48:ALA:HA	1:C:414:MET:HE2	1.87	0.55
1:D:348:GLY:HA2	1:D:353:ASN:HD22	1.71	0.55
1:B:50:ASN:HD21	1:B:421:SER:CB	2.19	0.55
1:C:2:LEU:HD21	1:C:36:LEU:CD2	2.36	0.55
1:C:295:MET:HE3	1:C:410:LYS:HD3	1.88	0.55
1:A:3:THR:O	1:A:7:GLU:HG3	2.07	0.54
1:A:80:GLY:HA3	1:B:314:THR:O	2.07	0.54
1:B:16:PRO:O	1:B:19:GLU:HG2	2.07	0.54
1:D:2:LEU:HD22	1:D:6:ARG:HD2	1.88	0.54
1:C:50:ASN:HD21	1:C:421:SER:CB	2.20	0.54
1:C:247:HIS:O	1:C:251:GLU:HG3	2.07	0.54
1:C:270:PRO:C	1:C:272:ASP:N	2.65	0.54
1:D:202:ALA:HA	1:D:228:ASP:OD1	2.07	0.54
1:C:394:LEU:HD13	1:C:411:THR:CG2	2.38	0.54
1:C:122:LEU:HD23	1:C:250:GLU:HB2	1.90	0.54
1:A:11:ALA:O	1:A:14:GLU:HB2	2.07	0.53
1:B:16:PRO:HA	1:B:19:GLU:HG2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ALA:HA	1:A:228:ASP:OD1	2.09	0.53
1:B:5:LEU:O	1:B:8:ASP:HB2	2.09	0.53
1:B:186:ALA:HB1	1:B:248:LEU:HD21	1.90	0.53
1:C:54:PHE:HB3	1:C:331:LEU:HD11	1.90	0.53
1:B:2:LEU:C	1:B:4:GLU:H	2.17	0.52
1:A:50:ASN:ND2	1:A:421:SER:OG	2.43	0.52
1:A:323:GLY:O	1:A:327:VAL:HG23	2.09	0.52
1:B:71:THR:HG21	1:B:98:LEU:HD11	1.90	0.52
1:B:115:VAL:HG21	1:B:279:LEU:HD13	1.90	0.52
1:A:30:THR:HB	1:A:430:GLU:HG3	1.89	0.52
1:D:184:LEU:HD11	1:D:193:LEU:HD13	1.92	0.52
1:D:91:VAL:HG22	1:D:125:ASP:HB3	1.91	0.52
1:A:270:PRO:C	1:A:272:ASP:N	2.67	0.52
1:C:351:ARG:NH1	1:C:415:SER:HB2	2.24	0.52
1:B:343:LEU:HD12	1:B:343:LEU:H	1.74	0.51
1:D:150:GLN:HE21	1:D:241:PRO:HD3	1.75	0.51
1:B:63:LEU:O	1:B:67:HIS:HB2	2.09	0.51
1:C:91:VAL:HG22	1:C:125:ASP:HB3	1.91	0.51
1:C:23:SER:HA	1:C:436:VAL:O	2.10	0.51
1:A:47:PRO:HB3	1:A:355:VAL:HG21	1.92	0.51
1:D:270:PRO:O	1:D:271:ALA:CB	2.59	0.51
1:B:133:ARG:HB3	1:B:150:GLN:HB3	1.92	0.50
1:A:54:PHE:CD2	1:A:363:LEU:HD22	2.46	0.50
1:C:51:MET:HE3	1:C:353:ASN:OD1	2.12	0.50
1:D:175:ALA:HB3	1:D:178:GLU:CD	2.36	0.50
1:B:213:PRO:HD2	1:B:218:THR:O	2.11	0.50
1:B:417:VAL:HG13	1:B:438:ASN:HD21	1.77	0.50
1:C:156:VAL:HG21	1:C:248:LEU:HD12	1.94	0.50
1:D:321:ASP:O	1:D:325:VAL:HG23	2.11	0.50
1:A:349:LEU:O	1:A:349:LEU:HD23	2.12	0.50
1:B:78:ALA:HB1	1:B:79:PRO:HD2	1.94	0.50
1:C:270:PRO:O	1:C:271:ALA:HB3	2.12	0.50
1:C:409:ALA:HA	1:C:422:GLY:HA3	1.94	0.50
1:A:58:ALA:HB2	1:A:327:VAL:HG13	1.94	0.50
1:B:30:THR:OG1	1:B:430:GLU:HG3	2.12	0.50
1:C:425:PRO:HA	1:C:430:GLU:OE2	2.12	0.50
1:D:386:SER:CB	1:D:399:ARG:NH2	2.75	0.50
1:D:397:ARG:HH12	1:D:449:GLN:NE2	2.09	0.50
1:B:361:VAL:HA	1:B:364:LEU:HD12	1.93	0.49
1:C:382:VAL:HB	1:C:385:GLU:HG3	1.94	0.49
1:D:361:VAL:HA	1:D:364:LEU:HD12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:HIS:O	1:A:250:GLU:HB3	2.12	0.49
1:D:44:GLN:HA	1:D:355:VAL:O	2.12	0.49
1:A:48:ALA:O	1:A:348:GLY:HA3	2.12	0.49
1:C:129:PHE:CB	1:C:153:ALA:HB2	2.42	0.49
1:D:8:ASP:O	1:D:12:ILE:HG13	2.12	0.49
1:A:184:LEU:HD11	1:A:193:LEU:HB2	1.95	0.49
1:D:25:VAL:HG22	1:D:435:ILE:HG23	1.94	0.49
1:B:304:GLU:OE1	1:B:304:GLU:HA	2.13	0.48
1:A:47:PRO:HB3	1:A:355:VAL:CG2	2.44	0.48
1:C:28:VAL:HG13	1:C:34:GLU:O	2.12	0.48
1:C:429:GLY:O	1:C:431:LEU:HD13	2.13	0.48
1:D:271:ALA:C	1:D:273:TRP:H	2.21	0.48
1:D:125:ASP:OD2	1:D:125:ASP:C	2.57	0.48
1:D:144:PRO:HG2	1:D:145:TYR:CE1	2.48	0.48
1:D:175:ALA:O	1:D:178:GLU:HG3	2.13	0.48
1:B:459:TYR:C	1:B:461:GLY:H	2.22	0.48
1:A:381:PRO:HD2	1:A:409:ALA:O	2.14	0.48
1:C:51:MET:CE	1:C:353:ASN:HB3	2.44	0.48
1:C:172:THR:HG22	1:C:231:PRO:HB3	1.96	0.48
1:C:147:TYR:HB2	1:C:300:ASN:ND2	2.29	0.48
1:D:28:VAL:HG12	1:D:29:ASP:N	2.28	0.48
1:D:133:ARG:HB3	1:D:150:GLN:HB3	1.96	0.48
1:A:50:ASN:HD21	1:A:421:SER:CB	2.27	0.47
1:D:250:GLU:HG2	1:D:260:VAL:HG21	1.96	0.47
1:A:47:PRO:CG	1:A:51:MET:HE2	2.39	0.47
1:B:270:PRO:HB2	1:B:273:TRP:CD1	2.49	0.47
1:C:295:MET:CE	1:C:410:LYS:HD3	2.45	0.47
1:A:28:VAL:HG13	1:A:34:GLU:O	2.14	0.47
1:C:46:LEU:HD11	1:C:351:ARG:O	2.15	0.47
1:C:321:ASP:O	1:C:325:VAL:HG23	2.14	0.47
1:D:88:LEU:HD21	1:D:279:LEU:HD12	1.97	0.47
1:C:32:THR:OG1	1:C:34:GLU:HB2	2.15	0.47
1:C:47:PRO:O	1:C:48:ALA:HB3	2.14	0.47
1:B:341:LEU:CD2	1:B:355:VAL:HG12	2.45	0.47
1:A:78:ALA:HB2	1:A:277:GLU:HG2	1.97	0.46
1:B:47:PRO:HB3	1:B:355:VAL:HG21	1.97	0.46
1:A:196:ARG:HB2	1:A:222:THR:HG22	1.97	0.46
1:A:366:GLN:HA	1:A:366:GLN:NE2	2.30	0.46
1:B:184:LEU:HD11	1:B:193:LEU:HB2	1.96	0.46
1:B:271:ALA:C	1:B:273:TRP:N	2.74	0.46
1:B:342:VAL:HB	1:B:354:LEU:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:MET:HE3	1:C:353:ASN:CG	2.41	0.46
1:A:292:VAL:HB	1:A:293:PRO:CD	2.46	0.46
1:C:11:ALA:O	1:C:14:GLU:HB2	2.16	0.46
1:C:46:LEU:HD13	1:C:353:ASN:O	2.16	0.46
1:A:88:LEU:HD21	1:A:279:LEU:HD12	1.97	0.46
1:B:331:LEU:O	1:B:334:LEU:HB3	2.16	0.46
1:C:133:ARG:HB3	1:C:150:GLN:HB3	1.97	0.46
1:A:139:TRP:HB3	1:A:141:GLU:OE1	2.16	0.46
1:B:47:PRO:O	1:B:50:ASN:HB2	2.15	0.46
1:C:129:PHE:CG	1:C:153:ALA:HB2	2.51	0.46
1:C:86:GLN:O	1:C:120:GLY:N	2.49	0.46
1:D:190:TYR:OH	1:D:243:ALA:HB3	2.15	0.46
1:B:25:VAL:HG13	1:B:435:ILE:HG12	1.98	0.46
1:C:54:PHE:CD2	1:C:363:LEU:HD22	2.51	0.45
1:D:106:MET:O	1:D:110:VAL:HG23	2.15	0.45
1:D:247:HIS:O	1:D:251:GLU:HG3	2.16	0.45
1:A:60:LEU:O	1:A:64:GLY:HA2	2.17	0.45
1:B:2:LEU:C	1:B:4:GLU:N	2.71	0.45
1:B:198:VAL:HG12	1:B:199:THR:N	2.32	0.45
1:B:270:PRO:C	1:B:272:ASP:H	2.22	0.45
1:C:101:GLU:N	1:C:101:GLU:OE1	2.49	0.45
1:D:211:ASP:OD1	1:D:212:ARG:N	2.49	0.45
1:B:122:LEU:HD23	1:B:250:GLU:HB2	1.98	0.45
1:C:78:ALA:HB2	1:C:277:GLU:OE2	2.17	0.45
1:B:345:ASP:OD1	1:B:345:ASP:C	2.60	0.45
1:A:125:ASP:C	1:A:125:ASP:OD2	2.59	0.45
1:D:51:MET:CE	1:D:353:ASN:HB3	2.45	0.45
1:D:10:ASP:O	1:D:14:GLU:HG2	2.17	0.45
1:B:128:TRP:CE2	1:B:312:GLN:HG3	2.52	0.44
1:B:299:ASN:C	1:B:299:ASN:ND2	2.65	0.44
1:C:253:LEU:HB3	1:C:258:VAL:HB	1.98	0.44
1:C:341:LEU:HD23	1:C:355:VAL:HA	1.98	0.44
1:D:381:PRO:HD2	1:D:409:ALA:O	2.17	0.44
1:B:37:TYR:CD1	1:B:38:SER:N	2.85	0.44
1:B:336:VAL:HG22	1:B:363:LEU:HA	1.99	0.44
1:C:49:SER:HB2	1:C:412:GLY:HA2	1.99	0.44
1:B:45:LEU:HB3	1:B:438:ASN:OD1	2.18	0.44
1:C:111:ALA:C	1:C:113:SER:H	2.25	0.44
1:D:44:GLN:HB3	1:D:354:LEU:HD22	1.98	0.44
1:A:119:ARG:NH1	1:A:261:LYS:HZ1	2.15	0.44
1:D:169:VAL:O	1:D:233:THR:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ARG:HE	1:C:119:ARG:HA	1.83	0.44
1:C:392:GLY:C	1:C:394:LEU:H	2.25	0.44
1:B:94:GLY:O	1:B:96:PRO:HD3	2.17	0.44
1:A:96:PRO:HB2	1:A:302:HIS:CD2	2.53	0.44
1:B:295:MET:HE3	1:B:410:LYS:HD3	1.99	0.44
1:B:372:TRP:O	1:B:374:GLN:N	2.50	0.44
1:C:424:VAL:HB	1:C:431:LEU:CB	2.47	0.44
1:D:5:LEU:HD22	1:D:459:TYR:CD1	2.53	0.44
1:A:417:VAL:HG13	1:A:438:ASN:HD21	1.82	0.44
1:C:78:ALA:HB1	1:C:79:PRO:HD2	2.00	0.44
1:D:60:LEU:HD21	1:D:291:LEU:HD11	2.00	0.44
1:D:342:VAL:HG21	1:D:354:LEU:HD12	2.00	0.44
1:B:325:VAL:O	1:B:329:GLU:HG3	2.18	0.43
1:B:464:ALA:HB3	4:B:2018:HOH:O	2.18	0.43
1:C:50:ASN:ND2	1:C:421:SER:OG	2.50	0.43
1:D:119:ARG:HD2	1:D:119:ARG:HA	1.83	0.43
1:D:150:GLN:NE2	1:D:241:PRO:HD3	2.32	0.43
1:B:341:LEU:HG	1:B:359:THR:HG21	2.01	0.43
1:C:299:ASN:ND2	1:C:302:HIS:H	2.16	0.43
1:D:142:ASP:HB3	1:D:147:TYR:CE1	2.54	0.43
1:A:65:ALA:HB1	1:A:288:SER:HB3	2.00	0.43
1:A:173:PRO:HG3	1:A:225:LEU:CD2	2.47	0.43
1:C:86:GLN:HA	1:C:119:ARG:HB3	2.00	0.43
1:A:121:ASP:HA	1:A:262:GLY:HA3	2.00	0.43
1:C:294:PHE:HE2	1:C:303:ALA:HB2	1.84	0.43
1:A:89:TYR:CD2	1:A:123:TYR:HB2	2.54	0.43
1:B:65:ALA:O	1:B:286:GLU:HB3	2.19	0.43
1:B:270:PRO:O	1:B:271:ALA:HB3	2.19	0.43
1:B:463:GLN:HE21	1:B:463:GLN:HB2	1.60	0.43
1:B:116:ARG:HB3	1:B:116:ARG:CZ	2.48	0.43
1:B:197:ALA:HB2	1:B:221:VAL:HG12	2.01	0.43
1:B:437:ASN:C	1:B:438:ASN:HD22	2.26	0.43
1:C:4:GLU:HG2	1:C:455:ARG:NH2	2.34	0.43
1:A:94:GLY:O	1:A:96:PRO:HD3	2.18	0.43
1:B:101:GLU:OE1	1:B:101:GLU:HA	2.18	0.43
1:D:386:SER:HA	1:D:399:ARG:HH21	1.84	0.43
1:C:63:LEU:O	1:C:67:HIS:HB2	2.19	0.42
1:C:299:ASN:C	1:C:299:ASN:HD22	2.27	0.42
1:D:49:SER:HB2	1:D:412:GLY:HA2	2.01	0.42
1:D:110:VAL:HG12	1:D:115:VAL:HB	2.00	0.42
1:B:372:TRP:O	1:B:373:ALA:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:GLY:O	1:C:262:GLY:HA3	2.20	0.42
1:B:41:GLY:HA2	1:B:357:ALA:HB3	2.00	0.42
1:C:1:ARG:HG2	1:C:1:ARG:HH11	1.83	0.42
1:C:96:PRO:HB2	1:C:302:HIS:CE1	2.54	0.42
1:B:408:GLU:O	1:B:422:GLY:HA3	2.19	0.42
1:D:29:ASP:OD1	1:D:32:THR:HG23	2.20	0.42
1:B:50:ASN:HD21	1:B:421:SER:HB2	1.83	0.42
1:D:47:PRO:O	1:D:48:ALA:HB3	2.19	0.42
1:D:79:PRO:HG3	1:D:115:VAL:HG22	2.01	0.42
1:D:253:LEU:HB3	1:D:258:VAL:HB	2.01	0.42
1:B:270:PRO:O	1:B:272:ASP:N	2.46	0.42
1:C:1:ARG:HG3	1:C:2:LEU:N	2.34	0.42
1:C:130:ASP:OD1	1:C:132:GLU:HB2	2.20	0.42
1:A:307:VAL:HG11	1:A:324:LEU:HD13	2.01	0.42
1:B:54:PHE:HD2	1:B:331:LEU:HD13	1.85	0.42
1:B:49:SER:HB2	1:B:412:GLY:HA2	2.01	0.42
1:B:273:TRP:O	1:B:274:GLN:C	2.63	0.42
1:C:45:LEU:HB2	1:C:436:VAL:HG11	2.02	0.42
1:D:232:VAL:HG12	1:D:233:THR:N	2.34	0.42
1:A:58:ALA:CB	1:A:327:VAL:HG13	2.50	0.42
1:B:142:ASP:HB3	1:B:147:TYR:CE1	2.55	0.42
1:B:89:TYR:CZ	1:B:270:PRO:HD3	2.55	0.41
1:C:190:TYR:OH	1:C:243:ALA:HB3	2.20	0.41
1:C:459:TYR:C	1:C:461:GLY:H	2.28	0.41
1:D:137:ASP:OD1	1:D:344:ASN:ND2	2.53	0.41
1:A:80:GLY:HA3	1:B:315:ALA:HA	2.02	0.41
1:B:119:ARG:HG2	1:B:119:ARG:HH11	1.85	0.41
1:C:226:PRO:O	1:C:229:ALA:HB3	2.20	0.41
1:A:312:GLN:HA	1:A:317:ALA:H	1.85	0.41
1:B:351:ARG:HG3	1:B:351:ARG:HH11	1.84	0.41
1:B:59:ALA:HA	1:B:310:ILE:HD11	2.02	0.41
1:C:401:THR:C	1:C:403:ALA:N	2.78	0.41
1:D:426:GLY:O	1:D:429:GLY:N	2.52	0.41
1:B:60:LEU:O	1:B:64:GLY:HA2	2.21	0.41
1:C:166:VAL:HG12	1:C:237:THR:HG22	2.02	0.41
1:C:430:GLU:O	1:C:431:LEU:HD12	2.19	0.41
1:C:304:GLU:OE1	1:C:304:GLU:HA	2.20	0.41
1:D:294:PHE:CD2	1:D:302:HIS:HB2	2.56	0.41
1:B:299:ASN:ND2	1:B:302:HIS:H	2.19	0.41
1:D:67:HIS:O	1:D:286:GLU:HA	2.21	0.41
1:D:256:ASN:N	1:D:256:ASN:ND2	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ALA:HA	1:A:277:GLU:O	2.21	0.41
1:A:141:GLU:CD	1:A:141:GLU:H	2.28	0.41
1:A:84:GLU:HG2	1:A:117:THR:HB	2.03	0.41
1:A:226:PRO:O	1:A:229:ALA:HB3	2.21	0.41
1:B:152:SER:O	1:B:305:MET:HE3	2.21	0.41
1:B:211:ASP:OD1	1:B:212:ARG:N	2.53	0.41
1:B:211:ASP:CG	1:B:212:ARG:N	2.79	0.41
1:B:383:ALA:HA	1:B:394:LEU:HB3	2.03	0.41
1:C:52:LYS:HE2	1:C:298:SER:C	2.46	0.41
1:C:294:PHE:CE2	1:C:303:ALA:HB2	2.55	0.41
1:C:454:VAL:O	1:C:458:GLU:HG3	2.21	0.41
1:D:100:ALA:O	1:D:103:LEU:HB2	2.21	0.41
1:B:116:ARG:NH2	1:B:116:ARG:CB	2.82	0.41
1:B:427:PRO:C	1:B:429:GLY:H	2.28	0.41
1:B:459:TYR:C	1:B:461:GLY:N	2.79	0.41
1:C:44:GLN:HA	1:C:355:VAL:O	2.21	0.41
1:B:66:ASP:O	1:B:286:GLU:HG2	2.20	0.40
1:C:2:LEU:HD11	1:C:34:GLU:OE2	2.22	0.40
1:A:110:VAL:O	1:A:113:SER:HB3	2.20	0.40
1:A:173:PRO:HG2	1:A:229:ALA:O	2.21	0.40
1:C:320:TRP:O	1:C:324:LEU:HD22	2.21	0.40
1:D:143:GLU:N	1:D:144:PRO:HD2	2.36	0.40
1:A:22:VAL:HG11	1:A:43:GLU:HG2	2.04	0.40
1:A:130:ASP:OD1	1:A:130:ASP:C	2.64	0.40
1:D:73:VAL:HA	1:D:89:TYR:O	2.21	0.40
1:B:190:TYR:OH	1:B:243:ALA:HB3	2.21	0.40
1:C:69:PHE:CD1	1:C:69:PHE:N	2.89	0.40
1:A:136:ASP:OD1	1:A:137:ASP:N	2.53	0.40
1:A:233:THR:HG22	1:A:234:ALA:N	2.35	0.40
1:B:91:VAL:HG22	1:B:125:ASP:CB	2.52	0.40
1:B:175:ALA:O	1:B:199:THR:HB	2.22	0.40
1:C:111:ALA:C	1:C:113:SER:N	2.80	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	465/489 (95%)	432 (93%)	29 (6%)	4 (1%)	14	28
1	B	464/489 (95%)	428 (92%)	34 (7%)	2 (0%)	30	48
1	C	464/489 (95%)	429 (92%)	32 (7%)	3 (1%)	21	38
1	D	465/489 (95%)	438 (94%)	26 (6%)	1 (0%)	43	62
All	All	1858/1956 (95%)	1727 (93%)	121 (6%)	10 (0%)	24	42

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	270	PRO
1	A	31	ALA
1	B	373	ALA
1	C	30	THR
1	A	81	ARG
1	D	86	GLN
1	A	339	ALA
1	B	86	GLN
1	C	373	ALA
1	C	465	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	339/356 (95%)	322 (95%)	17 (5%)	22	44
1	B	338/356 (95%)	325 (96%)	13 (4%)	29	54
1	C	338/356 (95%)	328 (97%)	10 (3%)	36	61
1	D	339/356 (95%)	324 (96%)	15 (4%)	25	49
All	All	1354/1424 (95%)	1299 (96%)	55 (4%)	27	52

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	GLU
1	A	50	ASN
1	A	88	LEU
1	A	98	LEU
1	A	141	GLU
1	A	154	LEU
1	A	211	ASP
1	A	250	GLU
1	A	270	PRO
1	A	272	ASP
1	A	299	ASN
1	A	312	GLN
1	A	324	LEU
1	A	341	LEU
1	A	349	LEU
1	A	455	ARG
1	A	463	GLN
1	B	18	LEU
1	B	43	GLU
1	B	50	ASN
1	B	88	LEU
1	B	98	LEU
1	B	154	LEU
1	B	198	VAL
1	B	299	ASN
1	B	324	LEU
1	B	328	GLU
1	B	349	LEU
1	B	374	GLN
1	B	388	PRO
1	C	19	GLU
1	C	43	GLU
1	C	50	ASN
1	C	154	LEU
1	C	170	SER
1	C	291	LEU
1	C	299	ASN
1	C	324	LEU
1	C	349	LEU
1	C	428	GLU
1	D	19	GLU
1	D	63	LEU

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Mol	Chain	Res	Type
1	D	88	LEU
1	D	154	LEU
1	D	170	SER
1	D	184	LEU
1	D	256	ASN
1	D	289	GLU
1	D	299	ASN
1	D	324	LEU
1	D	328	GLU
1	D	349	LEU
1	D	375	THR
1	D	376	TRP
1	D	413	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	50	ASN
1	A	256	ASN
1	A	299	ASN
1	A	312	GLN
1	A	366	GLN
1	A	396	ASN
1	A	449	GLN
1	A	463	GLN
1	B	50	ASN
1	B	150	GLN
1	B	256	ASN
1	B	299	ASN
1	B	344	ASN
1	B	366	GLN
1	B	396	ASN
1	B	437	ASN
1	B	449	GLN
1	B	463	GLN
1	C	50	ASN
1	C	256	ASN
1	C	299	ASN
1	C	344	ASN
1	C	366	GLN
1	C	437	ASN

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Mol	Chain	Res	Type
1	C	449	GLN
1	D	44	GLN
1	D	50	ASN
1	D	86	GLN
1	D	150	GLN
1	D	158	HIS
1	D	256	ASN
1	D	299	ASN
1	D	437	ASN
1	D	449	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 38 ligands modelled in this entry, 18 are monoatomic - leaving 20 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	C	500	-	4,4,4	0.36	0	6,6,6	0.14	0
2	SO4	D	502	-	4,4,4	0.40	0	6,6,6	0.11	0
2	SO4	A	504	-	4,4,4	0.44	0	6,6,6	0.10	0
2	SO4	D	500	-	4,4,4	0.34	0	6,6,6	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	C	503	-	4,4,4	0.37	0	6,6,6	0.10	0
2	SO4	B	503	-	4,4,4	0.42	0	6,6,6	0.15	0
2	SO4	A	500	-	4,4,4	0.35	0	6,6,6	0.13	0
2	SO4	C	501	-	4,4,4	0.40	0	6,6,6	0.10	0
2	SO4	B	500	-	4,4,4	0.37	0	6,6,6	0.13	0
2	SO4	C	504	-	4,4,4	0.43	0	6,6,6	0.09	0
2	SO4	A	501	-	4,4,4	0.35	0	6,6,6	0.07	0
2	SO4	D	501	-	4,4,4	0.34	0	6,6,6	0.13	0
2	SO4	A	502	-	4,4,4	0.37	0	6,6,6	0.14	0
2	SO4	B	501	-	4,4,4	0.37	0	6,6,6	0.11	0
2	SO4	C	502	-	4,4,4	0.37	0	6,6,6	0.16	0
2	SO4	B	502	-	4,4,4	0.41	0	6,6,6	0.08	0
2	SO4	B	504	-	4,4,4	0.40	0	6,6,6	0.09	0
2	SO4	A	503	-	4,4,4	0.37	0	6,6,6	0.10	0
2	SO4	D	503	-	4,4,4	0.36	0	6,6,6	0.08	0
2	SO4	D	504	-	4,4,4	0.42	0	6,6,6	0.09	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	467/489 (95%)	-0.48	2 (0%) 88 87	20, 37, 59, 86	0
1	B	466/489 (95%)	-0.28	6 (1%) 75 68	22, 45, 70, 90	0
1	C	466/489 (95%)	-0.28	2 (0%) 88 87	20, 45, 74, 88	0
1	D	467/489 (95%)	-0.47	2 (0%) 88 87	16, 35, 62, 85	0
All	All	1866/1956 (95%)	-0.38	12 (0%) 85 82	16, 40, 68, 90	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	464	ALA	2.8
1	A	271	ALA	2.7
1	C	465	PRO	2.7
1	B	31	ALA	2.6
1	D	275	ASP	2.5
1	B	3	THR	2.4
1	C	271	ALA	2.3
1	D	271	ALA	2.3
1	B	428	GLU	2.2
1	B	427	PRO	2.1
1	B	271	ALA	2.0
1	A	178	GLU	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 ⓘ

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
3	CO	C	513	1/1	0.29	0.17	119,119,119,119	0
3	CO	D	513	1/1	0.39	0.25	130,130,130,130	0
3	CO	A	513	1/1	0.55	0.19	108,108,108,108	0
3	CO	B	513	1/1	0.62	0.23	115,115,115,115	0
3	CO	C	512	1/1	0.76	0.24	116,116,116,116	0
3	CO	B	510	1/1	0.77	0.18	88,88,88,88	0
3	CO	C	510	1/1	0.78	0.18	85,85,85,85	0
3	CO	B	514	1/1	0.79	0.17	84,84,84,84	0
3	CO	C	514	1/1	0.80	0.23	79,79,79,79	0
2	SO4	C	504	5/5	0.80	0.12	85,86,87,89	0
3	CO	B	512	1/1	0.81	0.19	110,110,110,110	0
2	SO4	A	504	5/5	0.82	0.16	70,71,73,74	0
2	SO4	B	501	5/5	0.84	0.16	71,73,75,75	0
3	CO	A	512	1/1	0.85	0.15	93,93,93,93	0
2	SO4	C	502	5/5	0.86	0.12	77,77,79,80	0
3	CO	D	512	1/1	0.87	0.14	86,86,86,86	0
3	CO	A	514	1/1	0.87	0.28	83,83,83,83	0
3	CO	D	514	1/1	0.89	0.22	69,69,69,69	0
3	CO	D	511	1/1	0.91	0.13	49,49,49,49	0
3	CO	A	511	1/1	0.91	0.13	46,46,46,46	0
2	SO4	C	501	5/5	0.92	0.11	70,70,72,72	0
2	SO4	D	504	5/5	0.92	0.11	79,80,81,83	0
2	SO4	A	502	5/5	0.92	0.09	66,67,67,68	0
2	SO4	B	504	5/5	0.93	0.11	80,81,82,83	0
2	SO4	C	500	5/5	0.93	0.10	47,49,49,51	0
2	SO4	D	502	5/5	0.93	0.07	70,70,71,73	0
2	SO4	B	502	5/5	0.93	0.11	72,74,74,76	0
2	SO4	C	503	5/5	0.94	0.09	65,66,67,67	0
2	SO4	D	503	5/5	0.94	0.14	67,67,69,69	0
2	SO4	A	503	5/5	0.94	0.11	65,66,67,67	0
3	CO	A	510	1/1	0.94	0.11	54,54,54,54	0
2	SO4	B	503	5/5	0.95	0.10	59,59,59,61	0
2	SO4	D	501	5/5	0.95	0.09	63,64,65,65	0
2	SO4	A	501	5/5	0.97	0.08	60,60,61,62	0
2	SO4	B	500	5/5	0.98	0.06	51,53,53,54	0
3	CO	D	510	1/1	0.98	0.08	58,58,58,58	0
2	SO4	D	500	5/5	0.98	0.05	52,52,53,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	500	5/5	0.99	0.05	44,44,46,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.