



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 04:41 AM GMT

PDB ID : 2D3T
Title : Fatty Acid beta-oxidation multienzyme complex from Pseudomonas Fragi,
Form V
Authors : Tsuchiya, D.; Shimizu, N.; Ishikawa, M.; Suzuki, Y.; Morikawa, K.
Deposited on : 2005-10-01
Resolution : 3.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

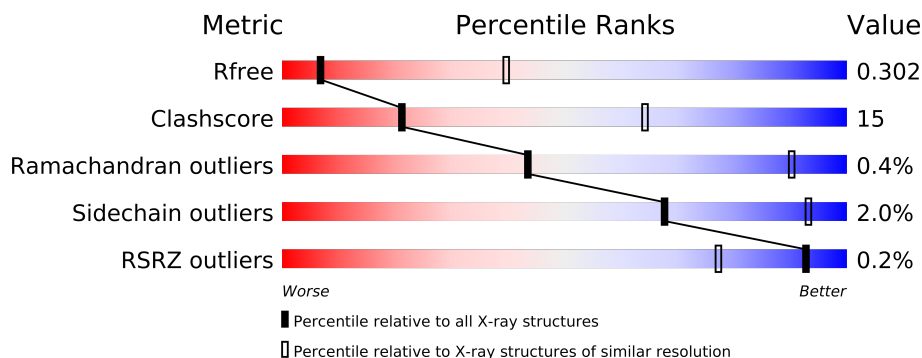
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.40 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}	66092	1017 (3.52-3.28)
Clashscore	79885	1214 (3.50-3.30)
Ramachandran outliers	78287	1177 (3.50-3.30)
Sidechain outliers	78261	1177 (3.50-3.30)
RSRZ outliers	66119	1017 (3.52-3.28)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	715	
1	B	715	
2	C	390	
2	D	390	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	ACO	C	1003	-	X
3	ACO	D	1004	-	X

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 16734 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

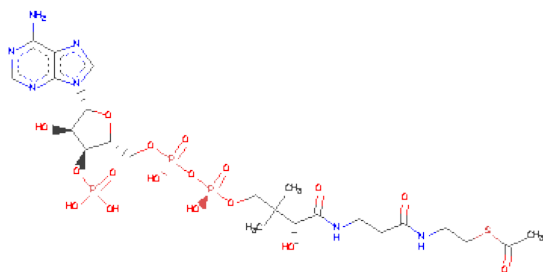
- Molecule 1 is a protein called Fatty oxidation complex alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	708	Total	C	N	O	S	0	0	0
			5368	3429	907	1005	27			
1	B	711	Total	C	N	O	S	0	0	0
			5390	3441	911	1011	27			

- Molecule 2 is a protein called 3-ketoacyl-CoA thiolase.

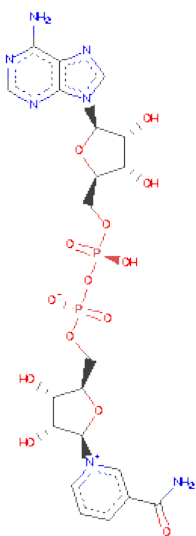
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	390	Total	C	N	O	S	0	0	0
			2893	1801	515	548	29			
2	D	390	Total	C	N	O	S	0	0	0
			2893	1801	515	548	29			

- Molecule 3 is ACETYL COENZYME *A (three-letter code: ACO) (formula: C₂₃H₃₈N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	C	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
3	D	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



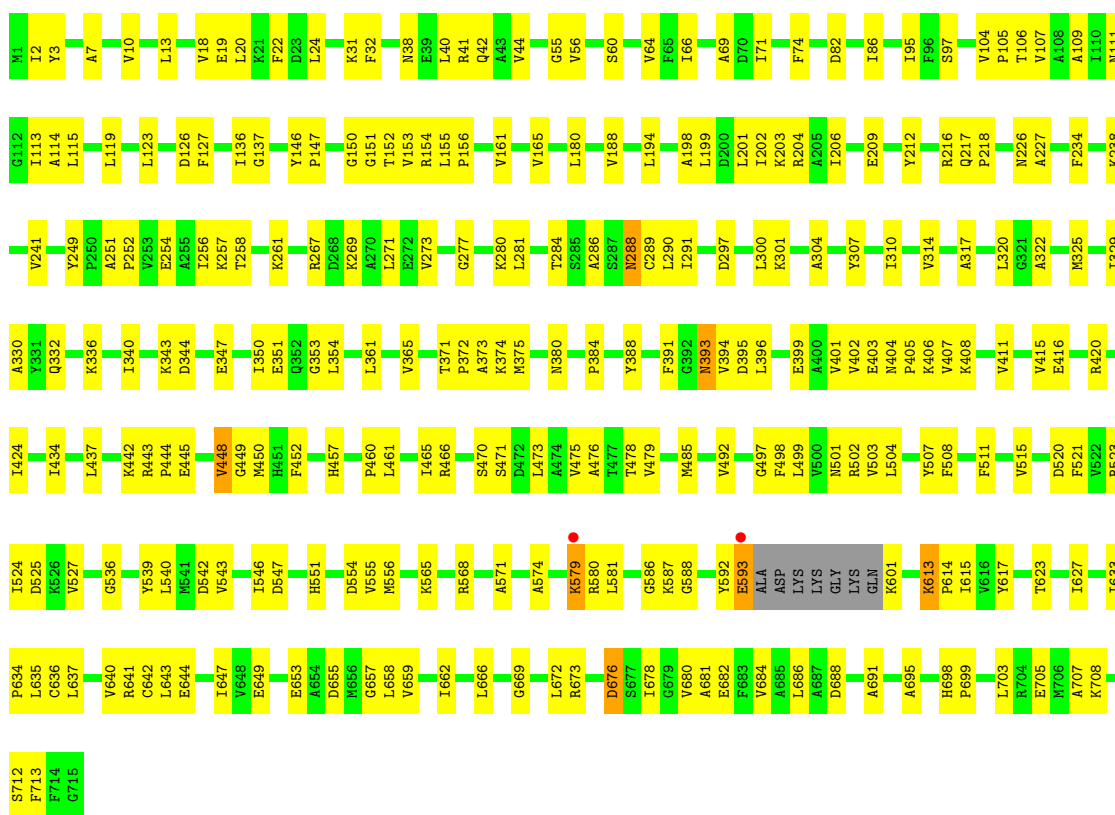
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
4	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

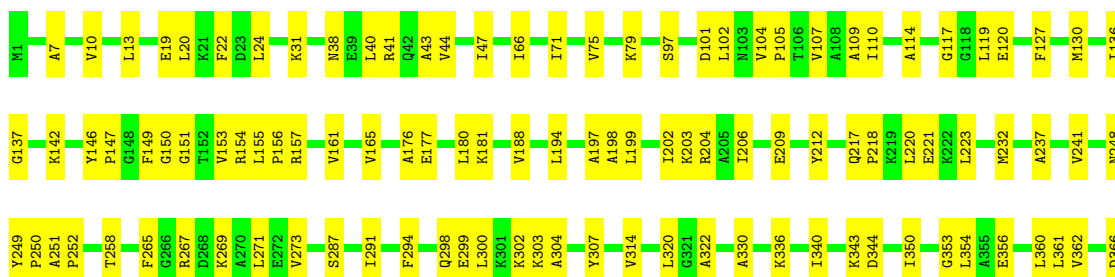
• Molecule 1: Fatty oxidation complex alpha subunit

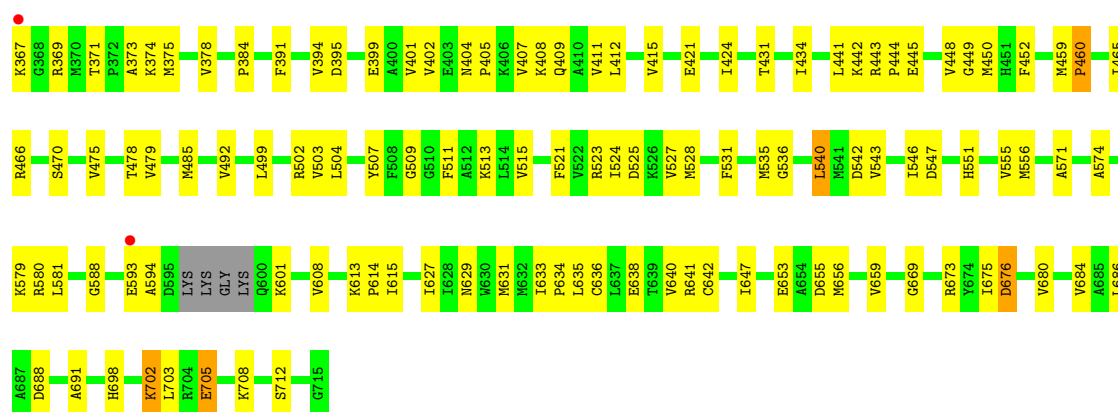
Chain A:



• Molecule 1: Fatty oxidation complex alpha subunit

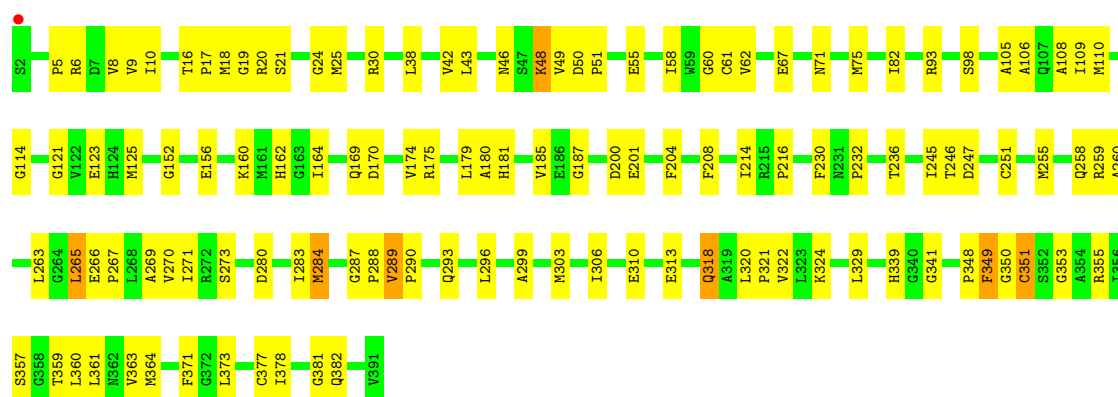
Chain B:





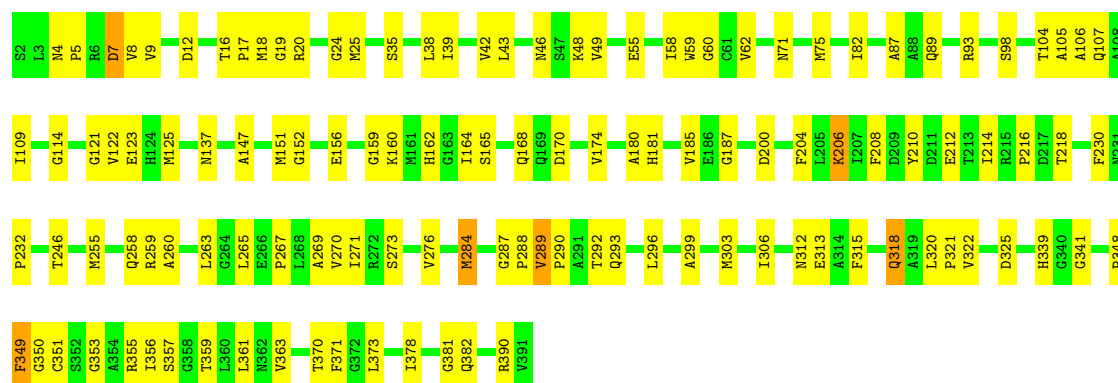
● Molecule 2: 3-ketoacyl-CoA thiolase

Chain C:



● Molecule 2: 3-ketoacyl-CoA thiolase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.23Å 137.55Å 198.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.40 13.79 – 3.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-3.40) 99.7 (13.79-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.29 (at 3.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.242 , 0.295 0.259 , 0.302	Depositor DCC
R_{free} test set	2509 reflections (7.00%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.843	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , -10.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 35837 reflections	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	16734	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.25	0/5454	0.46	0/7357
1	B	0.25	0/5476	0.46	0/7387
2	C	0.25	0/2941	0.47	0/3967
2	D	0.26	0/2941	0.47	0/3967
All	All	0.25	0/16812	0.46	0/22678

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5368	0	5494	179	0
1	B	5390	0	5511	156	0
2	C	2893	0	2903	96	0
2	D	2893	0	2903	97	0
3	C	51	0	34	0	0
3	D	51	0	34	2	0
4	A	44	0	26	0	0
4	B	44	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16734	0	16931	517	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (517) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:289:VAL:HG13	2:C:290:PRO:HD3	1.44	0.99
1:A:107:VAL:HG22	1:A:127:PHE:HB2	1.50	0.94
2:D:289:VAL:HG13	2:D:290:PRO:HD3	1.51	0.93
1:B:107:VAL:HG22	1:B:127:PHE:HB2	1.52	0.90
1:A:547:ASP:HB3	1:A:581:LEU:HD22	1.56	0.84
1:B:330:ALA:HA	1:B:340:ILE:HD13	1.60	0.83
1:A:330:ALA:HA	1:A:340:ILE:HD13	1.61	0.82
1:A:113:ILE:HD12	1:A:113:ILE:H	1.45	0.81
1:A:525:ASP:HB2	1:A:536:GLY:HA3	1.63	0.80
1:A:64:VAL:HG12	1:A:113:ILE:HD13	1.62	0.80
2:D:303:MET:HE3	2:D:306:ILE:HD12	1.63	0.80
1:A:371:THR:HB	1:A:374:LYS:HD3	1.63	0.79
2:C:306:ILE:HD13	2:C:373:LEU:HB2	1.65	0.79
1:A:109:ALA:HB1	1:A:194:LEU:HD21	1.66	0.78
2:C:263:LEU:HB2	2:C:265:LEU:HD22	1.66	0.77
1:A:372:PRO:HB2	1:B:362:VAL:HG13	1.66	0.76
2:D:62:VAL:HG21	2:D:349:PHE:HB3	1.68	0.76
1:B:547:ASP:HB3	1:B:581:LEU:HD22	1.68	0.75
1:A:669:GLY:HA3	1:A:673:ARG:HD2	1.69	0.74
1:B:525:ASP:HB2	1:B:536:GLY:HA3	1.70	0.74
2:D:125:MET:HE2	2:D:348:PRO:HA	1.69	0.74
1:A:640:VAL:HG11	1:A:703:LEU:HD12	1.70	0.73
1:A:71:ILE:HD11	1:A:291:ILE:HG23	1.67	0.73
2:C:62:VAL:HG21	2:C:349:PHE:HB3	1.71	0.72
1:B:71:ILE:HD11	1:B:291:ILE:HG23	1.71	0.71
1:A:373:ALA:HA	1:B:362:VAL:HG21	1.73	0.70
2:D:348:PRO:HB2	2:D:351:CYS:HB3	1.72	0.70
2:D:159:GLY:HA2	2:D:164:ILE:HD13	1.74	0.69
1:B:314:VAL:HG13	1:B:395:ASP:HB2	1.73	0.69
1:B:153:VAL:HG21	1:B:271:LEU:HD23	1.74	0.69
1:A:613:LYS:HB3	1:A:614:PRO:HD3	1.75	0.68
1:A:254:GLU:HG3	1:A:281:LEU:HD21	1.74	0.68
2:D:55:GLU:HG3	2:D:114:GLY:HA2	1.76	0.68
2:C:123:GLU:HB2	2:C:350:GLY:H	1.59	0.68
2:C:55:GLU:HG3	2:C:114:GLY:HA2	1.76	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:98:SER:HB3	2:D:353:GLY:HA3	1.77	0.67
2:D:287:GLY:O	2:D:290:PRO:HD2	1.95	0.67
2:D:265:LEU:HD23	2:D:265:LEU:H	1.60	0.67
2:C:287:GLY:O	2:C:290:PRO:HD2	1.95	0.66
1:A:465:ILE:H	1:A:465:ILE:HD12	1.59	0.66
2:D:306:ILE:HD13	2:D:373:LEU:HB2	1.76	0.66
1:A:251:ALA:HB3	1:A:252:PRO:HD3	1.78	0.66
2:D:378:ILE:HB	2:D:382:GLN:HB2	1.77	0.65
1:B:434:ILE:HD13	1:B:465:ILE:HG21	1.78	0.65
2:D:263:LEU:HB2	2:D:265:LEU:HD22	1.77	0.65
1:B:109:ALA:HB1	1:B:194:LEU:HD21	1.77	0.65
1:B:633:ILE:HB	1:B:634:PRO:HD3	1.79	0.65
1:B:251:ALA:HB3	1:B:252:PRO:HD3	1.78	0.64
2:C:125:MET:HE2	2:C:348:PRO:HA	1.78	0.64
1:B:180:LEU:HD12	1:B:188:VAL:HG23	1.78	0.64
1:B:613:LYS:HB3	1:B:614:PRO:HD3	1.79	0.64
1:A:354:LEU:HD11	1:A:384:PRO:HG2	1.79	0.64
2:C:289:VAL:HG23	2:C:293:GLN:HE21	1.63	0.64
1:A:642:CYS:HA	1:A:647:ILE:HD13	1.79	0.64
1:A:180:LEU:HD12	1:A:188:VAL:HG23	1.80	0.64
1:A:361:LEU:HB3	1:A:375:MET:HG3	1.80	0.63
2:C:265:LEU:HD23	2:C:265:LEU:H	1.64	0.63
1:A:633:ILE:HB	1:A:634:PRO:HD3	1.81	0.63
1:B:7:ALA:HA	1:B:24:LEU:HA	1.79	0.63
2:C:284:MET:HB3	2:C:381:GLY:H	1.64	0.63
1:A:637:LEU:HB3	1:A:641:ARG:HH12	1.62	0.63
2:C:348:PRO:HB2	2:C:351:CYS:HB3	1.80	0.63
2:D:16:THR:HG23	2:D:38:LEU:HD22	1.79	0.63
1:B:503:VAL:HG12	1:B:635:LEU:HG	1.80	0.62
2:D:25:MET:HE1	2:D:208:PHE:HB2	1.82	0.62
1:B:155:LEU:HB3	1:B:156:PRO:HD3	1.80	0.62
2:C:5:PRO:HB3	2:C:258:GLN:HB2	1.81	0.62
1:B:367:LYS:HB3	1:B:369:ARG:HE	1.65	0.62
2:C:156:GLU:O	2:C:160:LYS:HG2	1.99	0.62
1:A:7:ALA:HA	1:A:24:LEU:HA	1.82	0.61
1:B:336:LYS:HG3	1:B:485:MET:HA	1.81	0.61
2:D:359:THR:O	2:D:363:VAL:HG23	2.00	0.61
2:D:123:GLU:HB2	2:D:350:GLY:H	1.64	0.61
1:B:535:MET:SD	1:B:543:VAL:HG21	2.41	0.61
1:A:155:LEU:HB3	1:A:156:PRO:HD3	1.82	0.61
2:C:48:LYS:H	2:C:48:LYS:HD2	1.65	0.61
1:B:640:VAL:HG11	1:B:703:LEU:HD12	1.83	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:48:LYS:H	2:D:48:LYS:HD2	1.64	0.60
2:C:43:LEU:HD11	2:C:82:ILE:HD11	1.83	0.60
1:A:288:ASN:HD22	1:A:289:CYS:N	1.98	0.60
1:A:350:ILE:HG21	1:A:384:PRO:HB3	1.83	0.60
2:C:313:GLU:CD	2:C:341:GLY:HA3	2.21	0.60
1:B:705:GLU:HA	1:B:708:LYS:HE2	1.83	0.60
1:A:320:LEU:HD11	1:A:415:VAL:HG21	1.84	0.60
2:D:5:PRO:HB3	2:D:258:GLN:HB2	1.82	0.60
1:A:466:ARG:HD3	1:A:475:VAL:HG21	1.83	0.59
2:D:320:LEU:HB2	2:D:321:PRO:HD3	1.85	0.59
1:A:705:GLU:HA	1:A:708:LYS:HE2	1.84	0.59
1:A:226:ASN:ND2	1:A:227:ALA:H	2.01	0.59
1:A:307:TYR:HD2	1:A:476:ALA:HA	1.68	0.59
2:C:16:THR:HG23	2:C:38:LEU:HD22	1.84	0.59
2:D:273:SER:HB3	2:D:299:ALA:HB2	1.86	0.58
2:C:98:SER:HB3	2:C:353:GLY:HA3	1.86	0.58
2:D:180:ALA:HA	2:D:339:HIS:O	2.04	0.58
2:D:315:PHE:O	2:D:318:GLN:HG3	2.03	0.58
1:A:393:ASN:HD22	1:A:393:ASN:H	1.52	0.58
1:B:673:ARG:HH11	1:B:673:ARG:HB2	1.69	0.58
1:B:594:ALA:H	1:B:601:LYS:HD3	1.68	0.58
1:A:465:ILE:HA	1:A:492:VAL:O	2.03	0.57
2:C:25:MET:HE1	2:C:208:PHE:HB2	1.86	0.57
2:D:162:HIS:HB2	2:D:164:ILE:HD11	1.85	0.57
1:B:465:ILE:HA	1:B:492:VAL:O	2.04	0.57
2:D:9:VAL:HG13	2:D:267:PRO:HB3	1.87	0.57
2:C:320:LEU:HB2	2:C:321:PRO:HD3	1.87	0.57
1:B:350:ILE:HG21	1:B:384:PRO:HB3	1.86	0.57
1:A:404:ASN:H	1:A:408:LYS:HE2	1.70	0.57
1:B:300:LEU:HD11	1:B:656:MET:HG3	1.86	0.56
2:D:292:THR:O	2:D:296:LEU:HG	2.04	0.56
2:C:289:VAL:CG1	2:C:290:PRO:HD3	2.29	0.56
2:C:273:SER:HB3	2:C:299:ALA:HB2	1.87	0.56
1:B:343:LYS:HD3	1:B:344:ASP:N	2.20	0.56
2:C:8:VAL:HG11	2:C:271:ILE:HD12	1.86	0.56
1:B:7:ALA:HB2	1:B:24:LEU:HD13	1.88	0.56
2:C:162:HIS:HB2	2:C:164:ILE:HD11	1.87	0.56
2:C:9:VAL:HG13	2:C:267:PRO:HB3	1.88	0.56
1:A:695:ALA:HA	1:A:698:HIS:HD2	1.70	0.56
1:B:593:GLU:HB2	1:B:601:LYS:NZ	2.21	0.56
1:B:655:ASP:O	1:B:659:VAL:HG23	2.06	0.56
1:A:655:ASP:O	1:A:659:VAL:HG23	2.05	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:324:LYS:HB2	2:C:329:LEU:HD12	1.89	0.55
1:A:680:VAL:HG21	1:A:712:SER:HA	1.87	0.55
1:B:399:GLU:HG2	1:B:401:VAL:HG23	1.87	0.55
1:B:198:ALA:O	1:B:202:ILE:HG12	2.05	0.55
2:C:60:GLY:H	2:C:121:GLY:HA2	1.72	0.55
1:A:434:ILE:HD13	1:A:465:ILE:HG21	1.88	0.55
1:B:411:VAL:O	1:B:415:VAL:HG23	2.07	0.55
2:D:289:VAL:HG23	2:D:293:GLN:HE21	1.72	0.55
1:A:161:VAL:O	1:A:165:VAL:HG23	2.07	0.55
2:D:151:MET:HE3	3:D:1004:ACO:H21	1.89	0.55
1:B:442:LYS:HG3	1:B:443:ARG:HG3	1.89	0.55
1:A:399:GLU:HG2	1:A:401:VAL:HG23	1.87	0.55
1:B:161:VAL:O	1:B:165:VAL:HG23	2.07	0.55
1:A:521:PHE:HA	1:A:524:ILE:HD12	1.87	0.55
1:B:465:ILE:HD12	1:B:465:ILE:H	1.73	0.54
2:C:98:SER:HB2	2:C:350:GLY:O	2.07	0.54
1:A:314:VAL:HG13	1:A:395:ASP:HB2	1.90	0.54
1:B:102:LEU:HG	1:B:104:VAL:HG12	1.88	0.54
1:B:475:VAL:O	1:B:479:VAL:HG23	2.08	0.54
1:B:101:ASP:CG	1:B:267:ARG:HH22	2.11	0.54
1:A:258:THR:HG23	1:A:273:VAL:HG12	1.89	0.54
1:B:636:CYS:O	1:B:640:VAL:HG23	2.08	0.54
1:B:421:GLU:HA	1:B:443:ARG:NH2	2.23	0.54
2:D:98:SER:HB2	2:D:350:GLY:O	2.08	0.53
1:A:388:TYR:HE2	1:A:415:VAL:HG22	1.73	0.53
2:D:339:HIS:CD2	2:D:363:VAL:HG22	2.43	0.53
1:B:322:ALA:HB1	1:B:353:GLY:HA3	1.89	0.53
1:B:119:LEU:HD12	1:B:137:GLY:N	2.24	0.53
2:C:303:MET:HE3	2:C:306:ILE:HD12	1.91	0.53
2:D:98:SER:HB3	2:D:353:GLY:CA	2.38	0.53
1:A:411:VAL:O	1:A:415:VAL:HG23	2.09	0.53
1:A:461:LEU:HD22	1:A:657:GLY:O	2.09	0.53
2:C:378:ILE:HB	2:C:382:GLN:HB2	1.91	0.53
2:D:339:HIS:HD2	2:D:363:VAL:HG22	1.74	0.53
1:A:640:VAL:O	1:A:644:GLU:HG3	2.08	0.53
1:A:322:ALA:HB1	1:A:353:GLY:HA3	1.90	0.53
2:D:60:GLY:H	2:D:121:GLY:HA2	1.73	0.53
2:D:58:ILE:HD12	2:D:105:ALA:HB2	1.91	0.53
1:B:220:LEU:HD21	1:B:267:ARG:HH21	1.74	0.53
1:A:680:VAL:O	1:A:684:VAL:HG23	2.08	0.53
2:C:359:THR:O	2:C:363:VAL:HG23	2.09	0.53
1:B:507:TYR:CE1	1:B:631:MET:HB3	2.44	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:43:LEU:HD11	2:D:82:ILE:HD11	1.91	0.52
2:C:313:GLU:OE1	2:C:341:GLY:HA3	2.08	0.52
2:C:318:GLN:O	2:C:322:VAL:HG23	2.10	0.52
1:B:424:ILE:N	1:B:424:ILE:HD12	2.25	0.52
1:B:542:ASP:OD1	1:B:588:GLY:HA3	2.09	0.52
2:D:20:ARG:NH1	2:D:212:GLU:HB3	2.25	0.52
2:C:260:ALA:HA	2:C:265:LEU:HD21	1.92	0.52
1:B:571:ALA:HA	1:B:615:ILE:HG13	1.91	0.52
1:A:636:CYS:O	1:A:640:VAL:HG23	2.10	0.52
2:D:318:GLN:O	2:D:322:VAL:HG23	2.09	0.52
1:A:380:ASN:HB3	1:B:354:LEU:HD13	1.92	0.52
1:A:249:TYR:OH	1:A:666:LEU:HB2	2.10	0.51
2:C:180:ALA:HA	2:C:339:HIS:O	2.10	0.51
1:B:536:GLY:O	1:B:540:LEU:HB2	2.10	0.51
1:A:499:LEU:O	1:A:503:VAL:HG23	2.09	0.51
1:B:202:ILE:HG22	1:B:206:ILE:HD11	1.93	0.51
1:A:38:ASN:O	1:A:41:ARG:HB3	2.10	0.51
1:A:424:ILE:HD12	1:A:424:ILE:N	2.24	0.51
2:C:8:VAL:HG21	2:C:106:ALA:HB1	1.93	0.51
1:B:105:PRO:HG3	1:B:212:TYR:CD2	2.45	0.51
1:B:265:PHE:HB3	1:B:269:LYS:HB2	1.91	0.51
1:A:249:TYR:O	1:A:252:PRO:HD2	2.10	0.51
1:A:568:ARG:HD2	1:A:617:TYR:HE1	1.76	0.51
1:B:680:VAL:O	1:B:684:VAL:HG23	2.11	0.51
2:C:200:ASP:OD2	2:C:204:PHE:HB2	2.10	0.51
1:B:10:VAL:HG22	1:B:20:LEU:HG	1.93	0.51
1:A:204:ARG:O	1:A:209:GLU:HB3	2.11	0.51
2:C:10:ILE:HD11	2:C:271:ILE:HG13	1.93	0.50
2:D:8:VAL:HG11	2:D:271:ILE:HD12	1.92	0.50
1:A:150:GLY:HA3	1:A:154:ARG:NH1	2.26	0.50
2:D:306:ILE:HA	2:D:371:PHE:HB2	1.92	0.50
1:A:699:PRO:HB3	1:A:703:LEU:HD22	1.94	0.50
1:A:449:GLY:HA3	1:A:465:ILE:HB	1.93	0.50
2:C:38:LEU:O	2:C:42:VAL:HG23	2.12	0.50
1:B:523:ARG:O	1:B:527:VAL:HG23	2.11	0.50
1:B:551:HIS:O	1:B:555:VAL:HG23	2.11	0.50
1:A:336:LYS:HG3	1:A:485:MET:HA	1.94	0.50
2:C:105:ALA:O	2:C:109:ILE:HG13	2.11	0.50
2:C:318:GLN:C	2:C:321:PRO:HD2	2.31	0.50
2:D:62:VAL:CG2	2:D:349:PHE:HB3	2.40	0.50
2:D:164:ILE:N	2:D:164:ILE:HD12	2.26	0.50
2:D:16:THR:HG22	2:D:38:LEU:HD13	1.93	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:48:LYS:N	2:D:48:LYS:HD2	2.25	0.50
1:B:546:ILE:HD12	1:B:546:ILE:H	1.76	0.50
1:A:343:LYS:HD3	1:A:344:ASP:N	2.27	0.50
1:A:551:HIS:O	1:A:555:VAL:HG23	2.10	0.50
1:B:304:ALA:HA	1:B:307:TYR:HD1	1.76	0.50
1:B:492:VAL:HG13	1:B:647:ILE:CG2	2.42	0.50
1:A:571:ALA:HA	1:A:615:ILE:HG13	1.93	0.50
1:B:114:ALA:HB3	1:B:136:ILE:HG22	1.92	0.50
2:D:12:ASP:CB	2:D:46:ASN:HD21	2.24	0.50
1:A:217:GLN:HB3	1:A:218:PRO:HD3	1.93	0.49
1:A:405:PRO:HA	1:A:437:LEU:HD21	1.93	0.49
1:B:465:ILE:N	1:B:465:ILE:HD12	2.27	0.49
2:C:30:ARG:HD3	2:D:137:ASN:ND2	2.26	0.49
2:C:21:SER:OG	2:C:245:ILE:HG22	2.12	0.49
2:D:125:MET:HE3	2:D:246:THR:H	1.77	0.49
1:A:466:ARG:HG3	1:A:470:SER:HB2	1.95	0.49
1:B:269:LYS:O	1:B:273:VAL:HG23	2.12	0.49
2:C:21:SER:HA	2:C:247:ASP:OD2	2.12	0.49
1:A:105:PRO:HG3	1:A:212:TYR:CD2	2.48	0.49
1:B:204:ARG:O	1:B:209:GLU:HB3	2.13	0.49
1:B:367:LYS:HB3	1:B:369:ARG:NE	2.28	0.49
1:B:580:ARG:NH2	1:B:608:VAL:HG22	2.27	0.49
1:A:40:LEU:HD23	1:A:44:VAL:HG23	1.94	0.49
2:C:42:VAL:HG21	2:C:251:CYS:HB3	1.94	0.49
1:B:150:GLY:HA3	1:B:154:ARG:NH1	2.27	0.49
1:B:330:ALA:HA	1:B:340:ILE:HG21	1.94	0.49
2:D:255:MET:HE1	2:D:263:LEU:HD12	1.94	0.49
1:A:288:ASN:HD22	1:A:289:CYS:H	1.59	0.49
1:B:147:PRO:HB3	1:B:151:GLY:HA3	1.95	0.49
1:A:31:LYS:HB2	1:A:69:ALA:HA	1.95	0.49
1:A:119:LEU:HD12	1:A:137:GLY:N	2.27	0.49
2:C:164:ILE:HD12	2:C:164:ILE:N	2.27	0.49
2:D:123:GLU:HG2	2:D:349:PHE:HB2	1.95	0.49
2:C:93:ARG:HH12	2:D:87:ALA:HB1	1.77	0.49
1:A:542:ASP:OD1	1:A:588:GLY:HA3	2.13	0.48
2:D:8:VAL:O	2:D:270:VAL:HG13	2.14	0.48
1:A:393:ASN:HD22	1:A:393:ASN:N	2.10	0.48
1:A:676:ASP:OD2	1:A:712:SER:HB2	2.13	0.48
1:A:97:SER:HB3	1:A:267:ARG:NH1	2.28	0.48
1:B:638:GLU:HA	1:B:641:ARG:HH11	1.77	0.48
1:A:269:LYS:O	1:A:273:VAL:HG23	2.13	0.48
1:A:416:GLU:OE1	1:A:442:LYS:HG2	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:330:ALA:HA	1:A:340:ILE:HG21	1.95	0.48
1:A:40:LEU:O	1:A:44:VAL:HG23	2.13	0.48
1:A:146:TYR:HB2	1:A:147:PRO:HD2	1.96	0.48
1:B:199:LEU:HD11	1:B:203:LYS:HE3	1.94	0.48
1:A:288:ASN:HA	1:A:291:ILE:HD12	1.96	0.48
1:B:97:SER:HB3	1:B:267:ARG:NH1	2.29	0.48
2:D:259:ARG:NH2	2:D:263:LEU:HD21	2.28	0.48
2:D:260:ALA:HA	2:D:265:LEU:HD21	1.96	0.48
2:C:125:MET:CE	2:C:246:THR:H	2.27	0.48
1:A:508:PHE:HZ	1:A:540:LEU:HD22	1.77	0.48
1:B:371:THR:HG22	1:B:373:ALA:H	1.79	0.48
1:A:202:ILE:O	1:A:206:ILE:HG13	2.14	0.48
2:D:255:MET:HE3	2:D:259:ARG:HG3	1.95	0.47
1:A:202:ILE:HG22	1:A:206:ILE:HD11	1.96	0.47
2:C:280:ASP:HB2	2:C:283:ILE:HG12	1.96	0.47
1:A:22:PHE:CD2	1:A:66:ILE:HD11	2.49	0.47
2:D:206:LYS:HB2	2:D:206:LYS:NZ	2.28	0.47
1:A:504:LEU:O	1:A:507:TYR:HB3	2.14	0.47
1:A:127:PHE:CD2	1:A:201:LEU:HD21	2.49	0.47
1:B:354:LEU:HD21	1:B:384:PRO:HG3	1.95	0.47
1:B:521:PHE:HA	1:B:524:ILE:HD12	1.94	0.47
1:B:502:ARG:HA	1:B:556:MET:HE1	1.96	0.47
2:D:289:VAL:CG1	2:D:290:PRO:HD3	2.35	0.47
1:B:594:ALA:N	1:B:601:LYS:HD3	2.29	0.47
1:A:443:ARG:N	1:A:444:PRO:HD3	2.29	0.47
2:D:93:ARG:NH1	2:D:276:VAL:HG11	2.30	0.47
1:B:110:ILE:HB	1:B:130:MET:HG3	1.95	0.47
1:B:299:GLU:O	1:B:303:LYS:HG2	2.15	0.47
2:D:18:MET:HB3	2:D:246:THR:HG21	1.96	0.47
2:D:318:GLN:C	2:D:321:PRO:HD2	2.34	0.47
1:A:198:ALA:O	1:A:202:ILE:HG12	2.15	0.47
1:B:504:LEU:O	1:B:507:TYR:HB3	2.15	0.47
2:D:214:ILE:O	2:D:216:PRO:HD3	2.15	0.47
1:A:113:ILE:HG22	1:A:115:LEU:HG	1.97	0.47
1:A:475:VAL:O	1:A:479:VAL:HG23	2.13	0.47
2:D:105:ALA:O	2:D:109:ILE:HG13	2.14	0.47
1:A:147:PRO:HB3	1:A:151:GLY:HA3	1.96	0.47
2:C:214:ILE:O	2:C:216:PRO:HD3	2.15	0.47
1:B:248:ASN:O	1:B:250:PRO:HD3	2.14	0.47
1:A:593:GLU:H	1:A:593:GLU:CD	2.19	0.47
1:B:202:ILE:O	1:B:206:ILE:HG13	2.16	0.46
1:B:13:LEU:HD11	1:B:19:GLU:HB2	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:678:ILE:HG21	1:A:686:LEU:HD11	1.96	0.46
2:C:288:PRO:HB2	2:C:322:VAL:HG11	1.97	0.46
2:D:59:TRP:HE1	2:D:122:VAL:CG1	2.28	0.46
1:A:691:ALA:HB1	1:A:698:HIS:CE1	2.50	0.46
1:A:60:SER:HB3	1:A:111:ASN:OD1	2.16	0.46
2:C:108:ALA:HA	2:D:107:GLN:OE1	2.16	0.46
2:D:218:THR:HG23	3:D:1004:ACO:H61A	1.80	0.46
1:A:150:GLY:HA3	1:A:154:ARG:HH12	1.80	0.46
1:B:691:ALA:HB1	1:B:698:HIS:NE2	2.30	0.46
1:A:393:ASN:HA	1:A:420:ARG:NH1	2.30	0.46
1:A:13:LEU:HD11	1:A:19:GLU:HB2	1.98	0.46
1:B:669:GLY:HA3	1:B:673:ARG:HD2	1.97	0.46
2:C:296:LEU:HD21	2:C:306:ILE:HD11	1.96	0.46
1:B:673:ARG:HB2	1:B:673:ARG:NH1	2.30	0.46
1:A:373:ALA:HB2	1:B:362:VAL:HG11	1.97	0.46
2:C:18:MET:HB3	2:C:246:THR:HG21	1.98	0.46
2:C:8:VAL:O	2:C:270:VAL:HG13	2.16	0.46
1:A:32:PHE:CE1	1:A:66:ILE:HG21	2.50	0.46
1:B:499:LEU:O	1:B:503:VAL:HG23	2.16	0.46
2:C:152:GLY:HA3	2:C:230:PHE:CE2	2.51	0.46
1:A:649:GLU:HB3	1:A:653:GLU:OE1	2.16	0.46
2:D:8:VAL:HG21	2:D:106:ALA:HB1	1.97	0.46
1:A:391:PHE:HA	1:A:394:VAL:HG23	1.98	0.46
1:A:406:LYS:HA	1:A:406:LYS:HE2	1.97	0.46
2:D:59:TRP:HE1	2:D:122:VAL:HG12	1.81	0.45
1:A:119:LEU:HD23	1:A:123:LEU:HG	1.97	0.45
2:D:200:ASP:OD2	2:D:204:PHE:HB2	2.15	0.45
1:A:2:ILE:HD12	1:A:18:VAL:HG21	1.97	0.45
1:A:407:VAL:O	1:A:411:VAL:HG23	2.15	0.45
2:D:58:ILE:HD13	2:D:104:THR:HB	1.98	0.45
2:C:30:ARG:NH1	2:C:67:GLU:HG2	2.32	0.45
1:B:75:VAL:HG12	1:B:79:LYS:HE3	1.98	0.45
1:A:647:ILE:N	1:A:647:ILE:HD12	2.31	0.45
2:D:284:MET:HB3	2:D:381:GLY:H	1.82	0.45
1:B:157:ARG:HG2	1:B:223:LEU:HD23	1.98	0.45
1:A:241:VAL:HG21	1:A:256:ILE:HD11	1.99	0.45
2:D:19:GLY:O	2:D:246:THR:HG23	2.16	0.45
1:B:434:ILE:HD12	1:B:434:ILE:N	2.31	0.45
1:B:594:ALA:O	1:B:601:LYS:HB3	2.16	0.45
1:A:523:ARG:O	1:A:527:VAL:HG23	2.16	0.45
2:D:35:SER:O	2:D:39:ILE:HG13	2.17	0.45
1:A:465:ILE:HG12	1:A:498:PHE:CE2	2.52	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:20:ARG:O	2:D:24:GLY:HA3	2.17	0.45
1:B:258:THR:HG23	1:B:273:VAL:HG12	1.99	0.45
2:C:46:ASN:O	2:C:49:VAL:HG22	2.15	0.45
1:A:681:ALA:HA	1:A:707:ALA:HB1	1.99	0.45
1:B:459:MET:HA	1:B:460:PRO:HD3	1.77	0.45
1:A:613:LYS:HA	1:A:613:LYS:HE2	1.98	0.45
1:A:445:GLU:CD	1:A:445:GLU:H	2.20	0.45
2:C:17:PRO:HB2	2:C:25:MET:CE	2.46	0.45
1:B:320:LEU:HD11	1:B:415:VAL:HG21	1.98	0.45
1:A:347:GLU:O	1:A:351:GLU:HG2	2.17	0.45
2:C:98:SER:HB3	2:C:353:GLY:CA	2.47	0.45
1:B:546:ILE:N	1:B:546:ILE:HD12	2.32	0.45
2:D:156:GLU:O	2:D:160:LYS:HG2	2.17	0.45
1:B:580:ARG:CZ	1:B:608:VAL:HG22	2.48	0.44
1:B:391:PHE:HA	1:B:394:VAL:HG23	1.99	0.44
2:C:10:ILE:HD12	2:C:10:ILE:N	2.31	0.44
1:B:443:ARG:N	1:B:444:PRO:HD3	2.31	0.44
2:C:361:LEU:HA	2:C:364:MET:HE2	1.99	0.44
1:B:703:LEU:O	1:B:703:LEU:HD23	2.17	0.44
1:A:257:LYS:O	1:A:261:LYS:HG3	2.18	0.44
2:D:313:GLU:CD	2:D:341:GLY:HA3	2.37	0.44
2:C:306:ILE:HA	2:C:371:PHE:HB2	1.99	0.44
2:C:58:ILE:HD12	2:C:105:ALA:HB2	1.98	0.44
2:C:20:ARG:O	2:C:24:GLY:HA3	2.16	0.44
1:A:497:GLY:O	1:A:501:ASN:HB2	2.18	0.44
1:B:401:VAL:HG12	1:B:402:VAL:N	2.32	0.44
1:B:431:THR:HB	1:B:555:VAL:HG21	1.99	0.44
1:A:574:ALA:HB2	1:A:615:ILE:HD11	1.99	0.44
1:A:546:ILE:N	1:A:546:ILE:HD12	2.32	0.44
2:D:170:ASP:O	2:D:174:VAL:HG23	2.18	0.44
2:D:38:LEU:O	2:D:42:VAL:HG23	2.16	0.44
2:D:4:ASN:HB2	2:D:7:ASP:OD1	2.17	0.44
2:C:303:MET:HE1	2:C:373:LEU:HD22	2.00	0.44
2:C:19:GLY:O	2:C:246:THR:HG23	2.18	0.44
1:A:361:LEU:O	1:A:365:VAL:HG23	2.18	0.44
2:C:269:ALA:CB	2:C:361:LEU:HD22	2.48	0.44
1:B:177:GLU:HG2	1:B:181:LYS:HE3	2.00	0.44
1:B:642:CYS:O	1:B:647:ILE:HB	2.17	0.44
1:B:371:THR:HB	1:B:374:LYS:HG3	2.00	0.44
1:B:142:LYS:HB3	1:B:294:PHE:CE2	2.53	0.44
2:C:71:ASN:O	2:C:75:MET:HG2	2.18	0.44
1:B:197:ALA:HA	2:C:204:PHE:CE2	2.53	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:361:LEU:HB3	1:B:375:MET:HG3	1.99	0.44
1:B:702:LYS:O	1:B:702:LYS:HD2	2.18	0.43
1:A:10:VAL:HG22	1:A:20:LEU:HG	1.99	0.43
1:B:40:LEU:O	1:B:44:VAL:HG23	2.18	0.43
2:C:288:PRO:HD3	2:C:377:CYS:HB3	2.00	0.43
2:C:93:ARG:NH1	2:D:87:ALA:HB1	2.33	0.43
1:A:623:THR:O	1:A:627:ILE:HG13	2.18	0.43
2:D:152:GLY:HA3	2:D:230:PHE:CE2	2.53	0.43
1:B:676:ASP:OD2	1:B:712:SER:HB2	2.19	0.43
2:D:125:MET:CE	2:D:246:THR:H	2.31	0.43
1:B:407:VAL:O	1:B:411:VAL:HG23	2.18	0.43
1:B:217:GLN:HG3	1:B:221:GLU:OE1	2.19	0.43
2:C:164:ILE:HD11	2:C:321:PRO:HG3	2.00	0.43
2:C:265:LEU:H	2:C:265:LEU:CD2	2.29	0.43
1:A:520:ASP:O	1:A:524:ILE:HG13	2.18	0.43
1:B:43:ALA:O	1:B:47:ILE:HG13	2.17	0.43
2:C:19:GLY:H	2:C:246:THR:CG2	2.31	0.43
1:A:226:ASN:HD22	1:A:227:ALA:H	1.66	0.43
2:C:93:ARG:CZ	2:D:89:GLN:HB3	2.48	0.43
1:A:147:PRO:HB2	1:A:152:THR:HG23	2.00	0.43
1:B:528:MET:O	1:B:531:PHE:HB3	2.18	0.43
1:B:117:GLY:HA2	1:B:120:GLU:OE1	2.19	0.43
1:B:405:PRO:O	1:B:409:GLN:HG3	2.18	0.43
1:A:226:ASN:ND2	1:A:227:ALA:N	2.66	0.43
1:B:574:ALA:HB2	1:B:615:ILE:HD11	2.00	0.43
1:B:146:TYR:HB2	1:B:147:PRO:HD2	2.01	0.43
1:B:374:LYS:O	1:B:378:VAL:HG23	2.19	0.43
1:B:412:LEU:HB3	1:B:441:LEU:HD21	2.01	0.43
1:B:511:PHE:O	1:B:515:VAL:HG23	2.19	0.43
2:C:181:HIS:O	2:C:185:VAL:HG23	2.18	0.43
1:A:113:ILE:CD1	1:A:113:ILE:H	2.22	0.43
1:B:249:TYR:O	1:B:252:PRO:HD2	2.19	0.43
1:B:343:LYS:HB2	1:B:391:PHE:HZ	1.84	0.43
2:C:266:GLU:HA	2:C:267:PRO:HD3	1.92	0.43
1:A:40:LEU:HD23	1:A:40:LEU:O	2.19	0.43
2:C:310:GLU:HG3	2:C:360:LEU:HB2	2.01	0.43
1:A:310:ILE:O	1:A:473:LEU:HG	2.18	0.43
1:A:580:ARG:NH1	1:A:587:LYS:HB3	2.33	0.43
1:A:317:ALA:O	1:A:340:ILE:HA	2.18	0.43
1:B:267:ARG:O	1:B:271:LEU:HG	2.19	0.43
2:C:339:HIS:CE1	2:C:363:VAL:HG22	2.54	0.43
2:D:20:ARG:CZ	2:D:212:GLU:HB3	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:523:ARG:NH2	1:B:627:ILE:HD11	2.34	0.43
2:D:269:ALA:CB	2:D:361:LEU:HD22	2.49	0.43
2:C:175:ARG:NH1	2:C:179:LEU:HD11	2.33	0.43
2:D:122:VAL:HG22	2:D:123:GLU:N	2.34	0.43
2:D:17:PRO:HB2	2:D:25:MET:CE	2.49	0.43
1:A:688:ASP:HA	1:A:691:ALA:HB2	2.01	0.43
1:B:466:ARG:HD3	1:B:475:VAL:HG21	2.00	0.43
1:A:332:GLN:OE1	1:A:457:HIS:HA	2.19	0.43
1:A:658:LEU:O	1:A:662:ILE:HG22	2.18	0.43
1:A:448:VAL:HG21	1:A:478:THR:OG1	2.19	0.43
1:A:297:ASP:OD2	1:A:301:LYS:HE3	2.19	0.43
2:D:165:SER:OG	2:D:168:GLN:HG3	2.19	0.43
1:A:153:VAL:HG21	1:A:271:LEU:HD23	2.01	0.43
1:A:40:LEU:HD22	1:A:95:ILE:HG22	2.01	0.42
1:A:3:TYR:HE1	1:A:42:GLN:HE21	1.66	0.42
1:A:55:GLY:HA2	1:A:104:VAL:HG22	2.01	0.42
1:A:234:PHE:O	1:A:238:LYS:HG3	2.19	0.42
2:D:19:GLY:H	2:D:246:THR:CG2	2.31	0.42
1:A:434:ILE:HD12	1:A:434:ILE:N	2.34	0.42
1:B:629:ASN:HB3	1:B:633:ILE:HD11	2.00	0.42
2:C:17:PRO:HB2	2:C:25:MET:HE1	2.01	0.42
1:B:287:SER:O	1:B:291:ILE:HG13	2.20	0.42
1:B:492:VAL:HG13	1:B:647:ILE:HG23	2.00	0.42
2:D:17:PRO:HB3	2:D:210:TYR:O	2.19	0.42
2:D:71:ASN:O	2:D:75:MET:HG2	2.19	0.42
1:A:56:VAL:HB	1:A:106:THR:HG22	2.01	0.42
1:B:450:MET:HG2	1:B:452:PHE:CE1	2.55	0.42
2:D:181:HIS:O	2:D:185:VAL:HG23	2.19	0.42
1:A:325:MET:O	1:A:329:ILE:HG13	2.20	0.42
1:A:515:VAL:HG11	1:A:571:ALA:HB2	2.02	0.42
2:D:46:ASN:O	2:D:49:VAL:HG22	2.20	0.42
1:A:280:LYS:HB2	1:A:280:LYS:NZ	2.34	0.42
1:A:502:ARG:HA	1:A:556:MET:CE	2.49	0.42
1:A:401:VAL:HG12	1:A:402:VAL:N	2.35	0.42
1:A:442:LYS:HG3	1:A:443:ARG:HG3	2.01	0.42
1:B:120:GLU:CD	1:B:149:PHE:HB2	2.40	0.42
1:B:404:ASN:H	1:B:408:LYS:HE2	1.84	0.42
2:C:287:GLY:N	2:C:288:PRO:CD	2.82	0.42
2:D:287:GLY:N	2:D:288:PRO:CD	2.83	0.42
2:C:125:MET:HE3	2:C:246:THR:H	1.83	0.42
1:A:126:ASP:OD1	1:A:216:ARG:HB2	2.20	0.42
2:D:370:THR:HA	2:D:390:ARG:HB2	2.00	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:130:MET:CE	1:B:176:ALA:HA	2.49	0.42
1:A:502:ARG:HD3	1:A:642:CYS:SG	2.60	0.42
1:A:402:VAL:HG12	1:A:403:GLU:N	2.35	0.42
1:A:300:LEU:HD21	1:A:653:GLU:HG2	2.02	0.42
2:C:255:MET:HE1	2:C:263:LEU:HD12	2.01	0.41
1:B:525:ASP:CB	1:B:536:GLY:HA3	2.45	0.41
1:A:40:LEU:HD22	1:A:95:ILE:CG2	2.50	0.41
1:B:466:ARG:HG3	1:B:470:SER:HB2	2.02	0.41
1:A:586:GLY:HA2	1:A:592:TYR:HB2	2.02	0.41
1:A:539:TYR:O	1:A:543:VAL:HG23	2.20	0.41
1:B:232:MET:HG2	2:D:147:ALA:HB3	2.02	0.41
1:A:288:ASN:N	1:A:288:ASN:ND2	2.67	0.41
2:C:48:LYS:HD2	2:C:48:LYS:N	2.33	0.41
1:A:7:ALA:HB2	1:A:24:LEU:HD13	2.02	0.41
1:A:672:LEU:HD22	1:A:713:PHE:CE2	2.56	0.41
1:A:350:ILE:CG2	1:A:384:PRO:HB3	2.47	0.41
1:B:300:LEU:CD2	1:B:653:GLU:HG2	2.50	0.41
1:A:450:MET:HG2	1:A:452:PHE:CE1	2.55	0.41
2:D:312:ASN:HB2	2:D:356:ILE:HD13	2.02	0.41
1:A:199:LEU:HD11	1:A:203:LYS:HE3	2.02	0.41
1:B:579:LYS:HD2	1:B:579:LYS:N	2.35	0.41
1:B:298:GLN:O	1:B:302:LYS:HG3	2.20	0.41
1:B:636:CYS:HB3	1:B:675:ILE:HD11	2.03	0.41
1:A:146:TYR:OH	1:A:277:GLY:HA3	2.21	0.41
1:B:217:GLN:HB3	1:B:218:PRO:HD3	2.03	0.41
1:A:82:ASP:O	1:A:86:ILE:HG13	2.21	0.41
1:A:579:LYS:HA	1:A:579:LYS:HE2	2.02	0.41
1:B:350:ILE:CG2	1:B:384:PRO:HB3	2.49	0.41
2:C:169:GLN:HE22	2:C:236:THR:HB	1.85	0.41
1:B:445:GLU:H	1:B:445:GLU:CD	2.23	0.41
2:C:106:ALA:O	2:C:110:MET:HG3	2.21	0.41
1:A:396:LEU:HA	1:A:424:ILE:O	2.21	0.41
1:A:511:PHE:O	1:A:515:VAL:HG23	2.21	0.41
1:B:204:ARG:HG2	1:B:204:ARG:HH11	1.86	0.41
2:D:356:ILE:HG13	2:D:357:SER:N	2.36	0.41
1:B:509:GLY:O	1:B:513:LYS:HG3	2.20	0.41
1:A:284:THR:C	1:A:286:ALA:H	2.24	0.41
2:C:170:ASP:O	2:C:174:VAL:HG23	2.20	0.41
1:A:286:ALA:O	1:A:290:LEU:HG	2.21	0.41
1:B:448:VAL:HG21	1:B:478:THR:OG1	2.21	0.41
1:B:356:GLU:O	1:B:360:LEU:HG	2.21	0.41
1:B:22:PHE:CD2	1:B:66:ILE:HD11	2.57	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:288:ASN:N	1:A:288:ASN:HD22	2.19	0.40
1:A:492:VAL:HG13	1:A:647:ILE:CG2	2.51	0.40
1:B:449:GLY:HA3	1:B:465:ILE:HB	2.03	0.40
2:C:284:MET:HB2	2:C:378:ILE:O	2.21	0.40
1:A:471:SER:O	1:A:475:VAL:HG23	2.20	0.40
2:D:315:PHE:H	2:D:318:GLN:CG	2.34	0.40
1:B:688:ASP:HA	1:B:691:ALA:HB2	2.03	0.40
1:A:301:LYS:O	1:A:304:ALA:HB3	2.20	0.40
1:A:503:VAL:CG1	1:A:635:LEU:HG	2.52	0.40
2:C:201:GLU:H	2:C:201:GLU:CD	2.25	0.40
1:B:237:ALA:O	1:B:241:VAL:HG23	2.22	0.40
2:D:259:ARG:HH21	2:D:263:LEU:HD21	1.87	0.40
2:C:50:ASP:HA	2:C:51:PRO:HD3	1.98	0.40
2:C:255:MET:HE3	2:C:259:ARG:HG3	2.04	0.40
1:A:502:ARG:HG3	1:A:556:MET:HE1	2.02	0.40
2:C:357:SER:O	2:C:361:LEU:HG	2.22	0.40
1:B:40:LEU:HD23	1:B:44:VAL:HG23	2.04	0.40
1:B:38:ASN:O	1:B:41:ARG:HB3	2.22	0.40
2:C:6:ARG:HE	2:C:6:ARG:HA	1.87	0.40
1:A:114:ALA:HB3	1:A:136:ILE:HG22	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	704/715 (98%)	645 (92%)	57 (8%)	2 (0%)	50	92
1	B	707/715 (99%)	649 (92%)	57 (8%)	1 (0%)	59	96
2	C	388/390 (100%)	355 (92%)	30 (8%)	3 (1%)	27	81
2	D	388/390 (100%)	353 (91%)	32 (8%)	3 (1%)	27	81
All	All	2187/2210 (99%)	2002 (92%)	176 (8%)	9 (0%)	43	90

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	232	PRO
2	C	232	PRO
2	C	349	PHE
2	D	349	PHE
2	C	187	GLY
2	D	187	GLY
1	A	448	VAL
1	A	460	PRO
1	B	460	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	557/562 (99%)	545 (98%)	12 (2%)	64	93
1	B	559/562 (100%)	552 (99%)	7 (1%)	80	95
2	C	307/307 (100%)	299 (97%)	8 (3%)	59	91
2	D	307/307 (100%)	300 (98%)	7 (2%)	63	92
All	All	1730/1738 (100%)	1696 (98%)	34 (2%)	68	93

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	PHE
1	A	288	ASN
1	A	393	ASN
1	A	554	ASP
1	A	565	LYS
1	A	579	LYS
1	A	593	GLU
1	A	601	LYS
1	A	613	LYS
1	A	643	LEU
1	A	676	ASP
1	A	682	GLU
1	B	31	LYS
1	B	366	ASP

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Mol	Chain	Res	Type
1	B	540	LEU
1	B	676	ASP
1	B	686	LEU
1	B	702	LYS
1	B	705	GLU
2	C	48	LYS
2	C	61	CYS
2	C	265	LEU
2	C	284	MET
2	C	289	VAL
2	C	318	GLN
2	C	351	CYS
2	C	355	ARG
2	D	7	ASP
2	D	206	LYS
2	D	284	MET
2	D	289	VAL
2	D	318	GLN
2	D	325	ASP
2	D	355	ARG

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

Mol	Chain	Res	Type
1	A	38	ASN
1	A	42	GLN
1	A	226	ASN
1	A	288	ASN
1	A	296	ASN
1	A	298	GLN
1	A	393	ASN
1	A	409	GLN
1	A	488	ASN
1	A	493	ASN
1	A	629	ASN
1	A	698	HIS
1	B	38	ASN
1	B	42	GLN
1	B	163	ASN
1	B	260	GLN
1	B	296	ASN
1	B	298	GLN

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Mol	Chain	Res	Type
1	B	409	GLN
2	C	4	ASN
2	C	46	ASN
2	C	197	GLN
2	C	231	ASN
2	C	293	GLN
2	C	318	GLN
2	D	4	ASN
2	D	46	ASN
2	D	162	HIS
2	D	197	GLN
2	D	293	GLN
2	D	339	HIS
2	D	362	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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
4	NAD	A	1001	-	48,48,48	1.28	4 (8%)	73,73,73	1.45	7 (9%)
4	NAD	B	1002	-	48,48,48	1.27	4 (8%)	73,73,73	1.47	6 (8%)
3	ACO	C	1003	-	53,53,53	0.94	1 (1%)	79,79,79	0.96	2 (2%)
3	ACO	D	1004	-	53,53,53	0.99	1 (1%)	79,79,79	0.94	2 (2%)

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	NAD	A	1001	-	-	0/30/62/62	0/3/5/5
4	NAD	B	1002	-	-	0/30/62/62	0/3/5/5
3	ACO	C	1003	-	-	2/51/67/67	0/1/3/3
3	ACO	D	1004	-	-	2/51/67/67	0/1/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1001	NAD	C3N-C7N	5.11	1.59	1.50
4	B	1002	NAD	C3N-C7N	4.97	1.59	1.50
3	D	1004	ACO	CBP-CAP	4.55	1.60	1.55
3	C	1003	ACO	CBP-CAP	3.88	1.59	1.55
4	B	1002	NAD	C4N-C3N	3.40	1.45	1.39
4	A	1001	NAD	C4N-C3N	3.34	1.45	1.39
4	B	1002	NAD	C6N-N1N	3.20	1.44	1.35
4	A	1001	NAD	C2N-N1N	3.20	1.39	1.35
4	B	1002	NAD	C2N-N1N	3.19	1.39	1.35
4	A	1001	NAD	C6N-N1N	3.10	1.44	1.35

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1001	NAD	O4D-C1D-N1N	5.64	113.72	107.95
4	B	1002	NAD	O4D-C1D-N1N	5.63	113.72	107.95
4	B	1002	NAD	C6N-C5N-C4N	5.38	127.99	119.44
4	B	1002	NAD	C5N-C4N-C3N	-5.24	113.52	120.32
4	A	1001	NAD	C5N-C4N-C3N	-5.12	113.67	120.32
4	A	1001	NAD	C6N-C5N-C4N	5.05	127.47	119.44
3	C	1003	ACO	C2P-S1P-C	4.75	120.68	101.38
3	D	1004	ACO	C2P-S1P-C	4.67	120.35	101.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1003	ACO	P2A-O3A-P1A	-3.72	120.79	131.68
3	D	1004	ACO	P2A-O3A-P1A	-3.58	121.17	131.68
4	B	1002	NAD	C5N-C6N-N1N	-3.46	114.60	120.43
4	A	1001	NAD	C5N-C6N-N1N	-3.25	114.96	120.43
4	A	1001	NAD	C2N-C3N-C4N	2.97	121.68	118.31
4	B	1002	NAD	C2N-C3N-C4N	2.94	121.64	118.31
4	A	1001	NAD	PN-O3-PA	-2.65	121.55	132.95
4	B	1002	NAD	PN-O3-PA	-2.58	121.85	132.95
4	A	1001	NAD	O4D-C1D-C2D	-2.03	103.67	106.77

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	1003	ACO	CH3-C-S1P-C2P
3	C	1003	ACO	O-C-S1P-C2P
3	D	1004	ACO	CH3-C-S1P-C2P
3	D	1004	ACO	O-C-S1P-C2P

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	708/715 (99%)	0.05	2 (0%) 91 73	1, 34, 104, 149	0
1	B	711/715 (99%)	-0.10	2 (0%) 91 73	1, 18, 79, 148	0
2	C	390/390 (100%)	-0.22	1 (0%) 91 73	1, 4, 49, 126	0
2	D	390/390 (100%)	-0.14	0 100 100	1, 11, 68, 109	0
All	All	2199/2210 (99%)	-0.08	5 (0%) 93 77	1, 17, 85, 149	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	367	LYS	3.2
1	A	579	LYS	2.3
1	A	593	GLU	2.2
1	B	593	GLU	2.0
2	C	2	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ACO	C	1003	51/51	0.35	3.68	134,134,134,134	0
3	ACO	D	1004	51/51	0.42	2.18	184,184,184,184	0
4	NAD	B	1002	44/44	0.33	1.19	56,56,56,56	0
4	NAD	A	1001	44/44	0.35	0.52	57,57,57,57	0

6.5 Other polymers ⓘ

There are no such residues in this entry.