



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:43 pm GMT

PDB ID : 3TDL  
Title : Structure of human serum albumin in complex with DAUDA  
Authors : Wang, Y.; Luo, Z.; Shi, X.; Wang, H.; Nie, L.  
Deposited on : 2011-08-11  
Resolution : 2.60 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

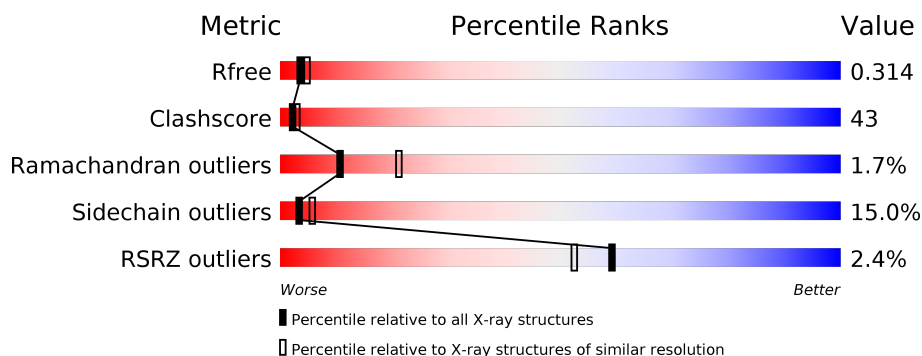
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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))
$R_{free}$	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	585	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MYR	A	1003	-	-	X	X
2	MYR	A	1004	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MYR	A	1005	-	-	-	X

## 2 Entry composition [i](#)

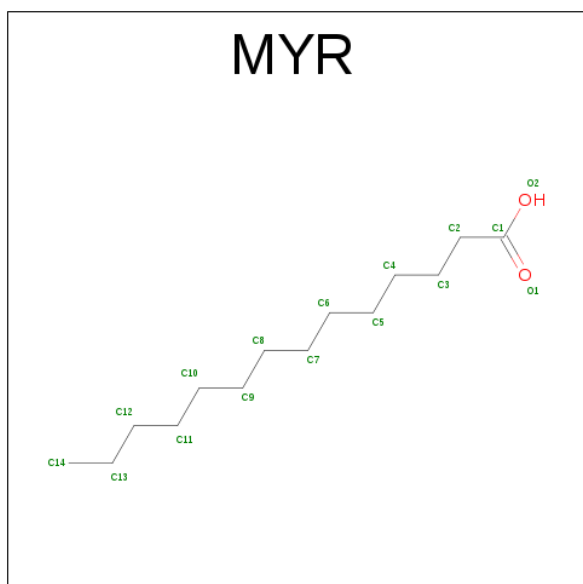
There are 3 unique types of molecules in this entry. The entry contains 4720 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Serum albumin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	578	Total	C	N	O	S	0	0	0
			4600	2906	778	875	41			

- Molecule 2 is MYRISTIC ACID (three-letter code: MYR) (formula:  $C_{14}H_{28}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	14	2		
2	A	1	Total	C	O	0	0
			16	14	2		
2	A	1	Total	C		0	0
			12	12			
2	A	1	Total	C	O	0	0
			16	14	2		

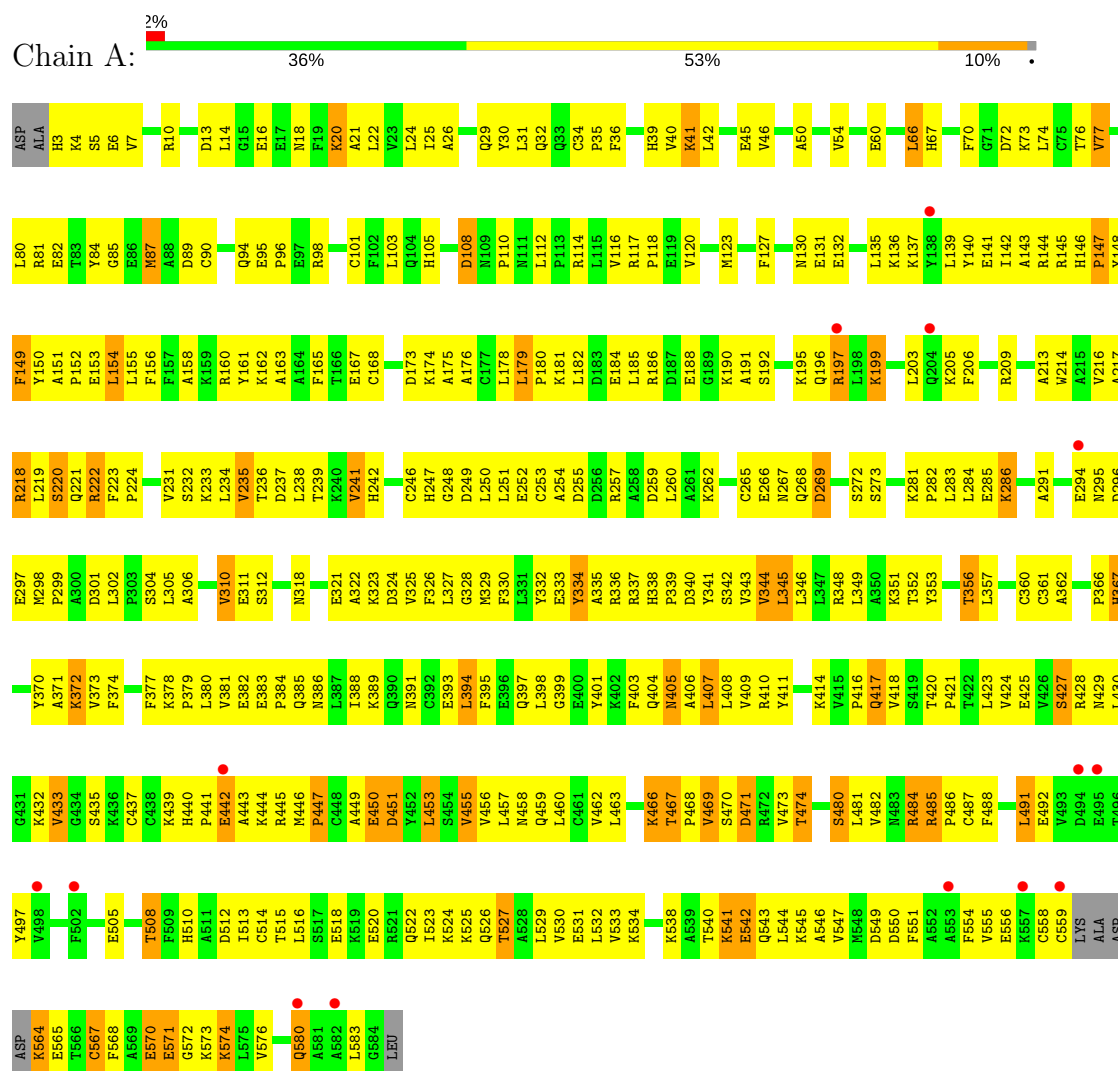
- Molecule 3 is 11-([5-(DIMETHYLAMINO)NAPHTHALEN-1-YL]SULFONYL)AMINO)U NDECANOIC ACID (three-letter code: 11D) (formula:  $C_{23}H_{34}N_2O_4S$ ).



### 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 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: Serum albumin



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	176.66Å 37.14Å 90.19Å 90.00° 104.38° 90.00°	Depositor
Resolution (Å)	35.50 – 2.60 35.50 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.2 (35.50-2.60) 97.7 (35.50-2.59)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.56 (at 2.58Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, $R_{free}$	0.251 , 0.261 0.251 , 0.314	Depositor DCC
$R_{free}$ test set	902 reflections (5.11%)	DCC
Wilson B-factor (Å <sup>2</sup> )	51.0	Xtriage
Anisotropy	0.632	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 34.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	4720	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.65% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 11D, MYR

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	0.49	0/4689	0.76	1/6323 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	PRO	N-CA-C	-5.74	97.17	112.10

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	4600	0	4525	381	0
2	A	60	0	104	41	0
3	A	60	0	66	16	0
All	All	4720	0	4695	403	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 43.

All (403) 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)
2:A:1003:MYR:H82	2:A:1003:MYR:C13	1.59	1.29
2:A:1004:MYR:C8	2:A:1004:MYR:H42	1.76	1.12
2:A:1003:MYR:H132	2:A:1003:MYR:H82	1.11	1.06
2:A:1004:MYR:C4	2:A:1004:MYR:H82	1.88	1.03
2:A:1004:MYR:H82	2:A:1004:MYR:H42	1.33	1.03
1:A:306:ALA:HA	1:A:310:VAL:HG23	1.40	1.03
1:A:523:ILE:O	1:A:527:THR:HG22	1.59	0.99
2:A:1003:MYR:H82	2:A:1003:MYR:C12	1.89	0.98
1:A:326:PHE:HE2	1:A:353:TYR:HH	1.04	0.97
1:A:385:GLN:O	1:A:388:ILE:HG13	1.65	0.95
1:A:467:THR:H	1:A:468:PRO:CD	1.70	0.95
1:A:467:THR:H	1:A:468:PRO:HD3	1.29	0.94
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.46	0.93
1:A:541:LYS:H	1:A:541:LYS:HD3	1.34	0.93
1:A:337:ARG:HH21	1:A:337:ARG:HG3	1.34	0.92
1:A:488:PHE:HB3	2:A:1004:MYR:H92	1.48	0.92
1:A:42:LEU:HD22	1:A:73:LYS:HE3	1.53	0.90
2:A:1004:MYR:C4	2:A:1004:MYR:C8	2.45	0.90
1:A:466:LYS:HA	1:A:466:LYS:HE3	1.53	0.90
1:A:467:THR:N	1:A:468:PRO:CD	2.35	0.89
2:A:1003:MYR:H132	2:A:1003:MYR:C8	2.00	0.86
1:A:16:GLU:O	1:A:20:LYS:HG3	1.77	0.85
1:A:31:LEU:HB2	1:A:39:HIS:HE1	1.39	0.85
1:A:342:SER:HB3	1:A:446:MET:HG2	1.57	0.84
1:A:410:ARG:O	1:A:414:LYS:HG3	1.77	0.84
1:A:81:ARG:NH1	1:A:89:ASP:OD2	2.11	0.83
1:A:259:ASP:O	1:A:262:LYS:HG2	1.79	0.82
1:A:236:THR:O	1:A:239:THR:HG22	1.79	0.82
1:A:467:THR:N	1:A:468:PRO:HD3	1.94	0.82
1:A:423:LEU:O	1:A:427:SER:HB2	1.78	0.81
1:A:505:GLU:HA	1:A:508:THR:HG22	1.63	0.80
1:A:341:TYR:CD1	1:A:345:LEU:HD13	2.17	0.79
1:A:249:ASP:OD1	1:A:252:GLU:HG2	1.84	0.78
1:A:353:TYR:HB2	1:A:377:PHE:HZ	1.46	0.78
2:A:1003:MYR:H122	2:A:1003:MYR:C8	2.14	0.78
1:A:108:ASP:O	1:A:110:PRO:HD3	1.84	0.77
1:A:31:LEU:HB2	1:A:39:HIS:CE1	2.19	0.77
1:A:222:ARG:HG2	1:A:222:ARG:HH21	1.47	0.76
1:A:442:GLU:HA	1:A:445:ARG:HD2	1.67	0.76
1:A:254:ALA:HB1	2:A:1002:MYR:H72	1.68	0.76
1:A:453:LEU:HD11	2:A:1004:MYR:C6	2.16	0.75
1:A:471:ASP:HA	1:A:474:THR:OG1	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:523:ILE:O	1:A:527:THR:CG2	2.34	0.75
1:A:152:PRO:O	1:A:155:LEU:HB3	1.87	0.75
1:A:22:LEU:HD13	2:A:1002:MYR:H82	1.67	0.75
1:A:3:HIS:HB2	1:A:6:GLU:OE2	1.88	0.74
2:A:1003:MYR:C12	2:A:1003:MYR:C8	2.58	0.73
1:A:424:VAL:O	1:A:428:ARG:HG3	1.89	0.72
1:A:428:ARG:HH21	1:A:526:GLN:HE22	1.37	0.72
1:A:306:ALA:HA	1:A:310:VAL:CG2	2.17	0.72
1:A:291:ALA:HB2	3:A:1007:11D:H24	1.71	0.72
1:A:394:LEU:HD23	1:A:397:GLN:HE21	1.54	0.72
1:A:233:LYS:NZ	1:A:237:ASP:OD1	2.23	0.72
1:A:254:ALA:CB	2:A:1002:MYR:H72	2.20	0.71
1:A:430:LEU:O	1:A:433:VAL:HG13	1.90	0.71
1:A:567:CYS:SG	1:A:568:PHE:N	2.65	0.70
1:A:45:GLU:OE2	1:A:73:LYS:NZ	2.21	0.70
1:A:282:PRO:HG2	1:A:285:GLU:HG3	1.74	0.69
1:A:168:CYS:HB2	1:A:178:LEU:HG	1.72	0.69
1:A:332:TYR:O	1:A:336:ARG:HG2	1.93	0.69
2:A:1004:MYR:H82	2:A:1004:MYR:H41	1.72	0.68
1:A:453:LEU:HD11	2:A:1004:MYR:H61	1.75	0.68
1:A:146:HIS:HB3	1:A:149:PHE:HB2	1.74	0.68
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.29	0.68
1:A:50:ALA:O	1:A:54:VAL:HG23	1.93	0.68
1:A:74:LEU:O	1:A:77:VAL:CG1	2.42	0.67
1:A:232:SER:HA	1:A:235:VAL:CG1	2.25	0.67
1:A:74:LEU:O	1:A:77:VAL:HG13	1.94	0.67
3:A:1006:11D:O16	3:A:1006:11D:H10	1.94	0.66
1:A:179:LEU:H	1:A:179:LEU:HD23	1.60	0.66
1:A:488:PHE:O	1:A:491:LEU:HB2	1.96	0.66
1:A:155:LEU:O	1:A:158:ALA:HB3	1.95	0.66
2:A:1003:MYR:H71	2:A:1004:MYR:H52	1.76	0.66
1:A:337:ARG:NH2	1:A:337:ARG:HG3	2.06	0.66
1:A:218:ARG:CG	1:A:343:VAL:HG21	2.27	0.65
1:A:397:GLN:HG3	1:A:398:LEU:HD13	1.78	0.65
1:A:576:VAL:HA	2:A:1005:MYR:H132	1.77	0.65
1:A:101:CYS:O	1:A:105:HIS:HD2	1.79	0.65
1:A:237:ASP:O	1:A:241:VAL:HG13	1.97	0.65
1:A:199:LYS:HB3	1:A:242:HIS:HE1	1.62	0.64
1:A:112:LEU:HD12	1:A:145:ARG:HG3	1.80	0.64
1:A:405:ASN:HD22	1:A:405:ASN:N	1.93	0.64
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:ARG:HG2	1:A:343:VAL:HG21	1.80	0.63
1:A:505:GLU:HA	1:A:508:THR:CG2	2.29	0.63
1:A:223:PHE:CD2	1:A:272:SER:HB2	2.33	0.63
1:A:418:VAL:HA	1:A:469:VAL:HG21	1.80	0.63
1:A:250:LEU:HD11	2:A:1002:MYR:H81	1.80	0.63
1:A:532:LEU:HD22	2:A:1005:MYR:H102	1.79	0.63
1:A:299:PRO:HD2	1:A:302:LEU:HD11	1.79	0.63
1:A:351:LYS:HG3	3:A:1006:11D:H27A	1.81	0.63
2:A:1003:MYR:C13	2:A:1004:MYR:H51	2.29	0.63
1:A:246:CYS:C	1:A:248:GLY:N	2.50	0.62
1:A:488:PHE:HA	1:A:491:LEU:HD22	1.81	0.62
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.82	0.62
1:A:216:VAL:HG23	1:A:235:VAL:HG11	1.82	0.62
1:A:429:ASN:O	1:A:433:VAL:HG12	1.99	0.62
1:A:84:TYR:HB3	1:A:87:MET:HB3	1.82	0.62
1:A:214:TRP:CZ3	3:A:1007:11D:H12A	2.35	0.61
1:A:120:VAL:HG11	1:A:174:LYS:HB3	1.81	0.61
1:A:525:LYS:HE2	1:A:551:PHE:HE2	1.64	0.61
1:A:395:PHE:HE2	1:A:439:LYS:HZ1	1.44	0.61
1:A:21:ALA:O	1:A:25:ILE:HG13	2.01	0.61
1:A:173:ASP:HB3	1:A:176:ALA:HB3	1.82	0.61
1:A:206:PHE:CD1	1:A:481:LEU:HD13	2.36	0.60
1:A:306:ALA:CA	1:A:310:VAL:HG23	2.22	0.60
1:A:485:ARG:HD3	2:A:1003:MYR:O1	2.01	0.60
2:A:1003:MYR:H141	2:A:1004:MYR:H51	1.83	0.60
1:A:156:PHE:CE1	1:A:285:GLU:HA	2.37	0.60
1:A:516:LEU:HB3	1:A:520:GLU:HB2	1.82	0.60
1:A:382:GLU:O	1:A:386:ASN:N	2.28	0.60
1:A:395:PHE:CZ	1:A:435:SER:HA	2.36	0.59
1:A:352:THR:O	1:A:356:THR:HG23	2.02	0.59
1:A:168:CYS:CB	1:A:178:LEU:HG	2.32	0.59
1:A:216:VAL:O	1:A:220:SER:OG	2.17	0.59
1:A:408:LEU:HG	1:A:530:VAL:HG22	1.83	0.59
2:A:1003:MYR:C13	2:A:1003:MYR:C8	2.56	0.59
1:A:353:TYR:HB2	1:A:377:PHE:CZ	2.35	0.59
1:A:150:TYR:HH	2:A:1002:MYR:C1	2.15	0.59
1:A:408:LEU:HG	1:A:530:VAL:CG2	2.32	0.58
1:A:541:LYS:HD3	1:A:541:LYS:N	2.14	0.58
1:A:394:LEU:HD22	1:A:398:LEU:HD22	1.85	0.58
1:A:337:ARG:HH21	1:A:337:ARG:CG	2.14	0.58
1:A:348:ARG:CG	1:A:482:VAL:HB	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ALA:HB2	3:A:1006:11D:H19	1.85	0.58
1:A:246:CYS:O	1:A:248:GLY:N	2.37	0.58
1:A:281:LYS:HB3	1:A:282:PRO:HD2	1.85	0.58
1:A:217:ALA:HB3	1:A:343:VAL:HG13	1.86	0.57
1:A:156:PHE:CD2	1:A:156:PHE:N	2.72	0.57
1:A:459:GLN:HG2	1:A:460:LEU:N	2.19	0.57
1:A:214:TRP:HZ3	3:A:1007:11D:H12A	1.69	0.57
1:A:342:SER:HB2	1:A:450:GLU:HB3	1.85	0.57
1:A:219:LEU:HD11	1:A:238:LEU:HD12	1.86	0.57
1:A:95:GLU:O	1:A:96:PRO:C	2.41	0.57
1:A:232:SER:HA	1:A:235:VAL:HG12	1.87	0.57
1:A:485:ARG:HH11	1:A:486:PRO:HA	1.70	0.57
1:A:547:VAL:O	1:A:550:ASP:HB2	2.04	0.57
1:A:395:PHE:CE1	1:A:435:SER:HA	2.39	0.56
1:A:526:GLN:O	1:A:530:VAL:HG23	2.05	0.56
1:A:291:ALA:HA	3:A:1007:11D:H20A	1.87	0.56
1:A:241:VAL:HG11	1:A:260:LEU:HD22	1.86	0.56
1:A:326:PHE:HE2	1:A:353:TYR:OH	1.82	0.56
1:A:132:GLU:O	1:A:136:LYS:HG2	2.05	0.56
1:A:156:PHE:O	1:A:160:ARG:HG3	2.06	0.56
1:A:341:TYR:CE1	1:A:381:VAL:HG22	2.41	0.56
1:A:160:ARG:HH12	1:A:188:GLU:CD	2.09	0.56
1:A:249:ASP:OD1	1:A:252:GLU:CG	2.54	0.56
1:A:444:LYS:O	1:A:447:PRO:HD2	2.05	0.56
1:A:332:TYR:CE1	1:A:336:ARG:HD2	2.41	0.55
1:A:153:GLU:OE2	1:A:192:SER:OG	2.23	0.55
1:A:246:CYS:C	1:A:248:GLY:H	2.08	0.55
1:A:395:PHE:CD2	1:A:395:PHE:C	2.79	0.55
1:A:334:TYR:CE1	1:A:338:HIS:HD2	2.24	0.55
1:A:449:ALA:O	1:A:453:LEU:HB2	2.06	0.55
1:A:471:ASP:HA	1:A:474:THR:HG1	1.71	0.55
1:A:150:TYR:OH	2:A:1002:MYR:C1	2.55	0.55
1:A:238:LEU:HD21	3:A:1007:11D:H21A	1.89	0.55
1:A:466:LYS:CA	1:A:468:PRO:HD3	2.37	0.54
1:A:567:CYS:O	1:A:571:GLU:N	2.39	0.54
1:A:332:TYR:CZ	1:A:336:ARG:HD2	2.42	0.54
1:A:337:ARG:NH2	1:A:337:ARG:CG	2.70	0.54
1:A:378:LYS:N	1:A:379:PRO:HD2	2.21	0.54
1:A:543:GLN:HB3	1:A:583:LEU:CD2	2.38	0.54
1:A:142:ILE:CG2	1:A:154:LEU:HD21	2.38	0.54
1:A:343:VAL:N	1:A:450:GLU:OE1	2.34	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:ARG:HD2	1:A:526:GLN:HE22	1.72	0.54
1:A:437:CYS:O	1:A:440:HIS:HB2	2.08	0.54
1:A:135:LEU:HD11	1:A:162:LYS:HB2	1.88	0.53
1:A:323:LYS:O	1:A:327:LEU:HG	2.09	0.53
1:A:381:VAL:O	1:A:384:PRO:HD2	2.08	0.53
1:A:411:TYR:CE1	2:A:1004:MYR:H112	2.43	0.53
1:A:455:VAL:HG12	1:A:456:VAL:N	2.23	0.53
1:A:338:HIS:HB3	1:A:341:TYR:CD2	2.43	0.53
1:A:430:LEU:HA	1:A:433:VAL:CG1	2.38	0.53
1:A:103:LEU:HD21	1:A:249:ASP:HB2	1.89	0.53
1:A:32:GLN:O	1:A:144:ARG:NH2	2.42	0.53
1:A:466:LYS:HA	1:A:466:LYS:CE	2.34	0.53
1:A:453:LEU:HD12	1:A:485:ARG:HG3	1.91	0.53
1:A:259:ASP:O	1:A:262:LYS:CG	2.55	0.53
1:A:480:SER:HG	1:A:482:VAL:CG2	2.22	0.53
1:A:351:LYS:HA	3:A:1006:11D:H23A	1.91	0.53
1:A:120:VAL:HG11	1:A:174:LYS:CB	2.39	0.53
1:A:340:ASP:HA	1:A:443:ALA:O	2.08	0.53
1:A:345:LEU:HD12	1:A:446:MET:SD	2.49	0.53
1:A:179:LEU:N	1:A:179:LEU:HD23	2.24	0.53
1:A:408:LEU:HD12	1:A:427:SER:CB	2.40	0.52
1:A:428:ARG:HH21	1:A:526:GLN:NE2	2.05	0.52
1:A:155:LEU:HD23	1:A:284:LEU:HD21	1.92	0.52
1:A:156:PHE:CZ	1:A:284:LEU:HG	2.44	0.52
1:A:179:LEU:N	1:A:179:LEU:CD2	2.72	0.52
1:A:266:GLU:C	1:A:267:ASN:HD22	2.13	0.52
1:A:420:THR:OG1	1:A:534:LYS:NZ	2.35	0.52
1:A:381:VAL:CG1	1:A:385:GLN:OE1	2.58	0.52
1:A:414:LYS:HA	1:A:492:GLU:O	2.08	0.52
1:A:167:GLU:OE1	1:A:181:LYS:HE3	2.10	0.52
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.30	0.52
2:A:1003:MYR:C14	2:A:1004:MYR:H51	2.41	0.51
1:A:485:ARG:NH2	2:A:1003:MYR:O1	2.41	0.51
1:A:253:CYS:O	1:A:257:ARG:HG3	2.11	0.51
1:A:218:ARG:HG3	1:A:343:VAL:HG21	1.93	0.51
1:A:367:HIS:O	1:A:371:ALA:HB2	2.11	0.51
1:A:411:TYR:HB3	2:A:1004:MYR:H142	1.93	0.51
1:A:380:LEU:O	1:A:384:PRO:CD	2.58	0.51
1:A:197:ARG:NH1	1:A:458:ASN:HD21	2.08	0.51
1:A:186:ARG:HD2	1:A:190:LYS:HD3	1.92	0.50
1:A:458:ASN:O	1:A:462:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ARG:CG	1:A:222:ARG:HH21	2.18	0.50
1:A:318:ASN:O	1:A:321:GLU:HG2	2.12	0.50
1:A:221:GLN:NE2	1:A:341:TYR:O	2.43	0.50
1:A:342:SER:HB2	1:A:450:GLU:CB	2.41	0.50
1:A:407:LEU:HG	1:A:411:TYR:HE2	1.74	0.50
1:A:525:LYS:HE2	1:A:551:PHE:CE2	2.46	0.50
1:A:394:LEU:O	1:A:397:GLN:HG2	2.12	0.50
1:A:218:ARG:HG2	1:A:343:VAL:CG2	2.41	0.50
1:A:424:VAL:HG11	1:A:523:ILE:HG23	1.93	0.50
1:A:147:PRO:O	1:A:148:TYR:HB2	2.12	0.50
1:A:405:ASN:O	1:A:409:VAL:HG13	2.12	0.50
1:A:518:GLU:O	1:A:522:GLN:HG3	2.12	0.50
1:A:529:LEU:O	1:A:532:LEU:HB3	2.12	0.50
1:A:334:TYR:CE1	1:A:338:HIS:CD2	3.00	0.50
1:A:318:ASN:HA	1:A:321:GLU:HG2	1.92	0.49
1:A:437:CYS:O	1:A:445:ARG:HG2	2.12	0.49
1:A:72:ASP:O	1:A:76:THR:HG23	2.12	0.49
1:A:39:HIS:CD2	1:A:140:TYR:HE2	2.30	0.49
1:A:524:LYS:HA	1:A:527:THR:HG23	1.94	0.49
1:A:137:LYS:O	1:A:141:GLU:HG2	2.12	0.49
1:A:36:PHE:CE1	1:A:137:LYS:HA	2.47	0.49
1:A:485:ARG:NH1	1:A:486:PRO:HA	2.27	0.49
1:A:540:THR:C	1:A:542:GLU:N	2.65	0.49
1:A:231:VAL:O	1:A:235:VAL:HG12	2.13	0.49
1:A:348:ARG:HG2	1:A:482:VAL:HB	1.94	0.49
1:A:220:SER:O	1:A:336:ARG:HA	2.13	0.49
1:A:186:ARG:HD2	1:A:190:LYS:CD	2.42	0.49
1:A:262:LYS:O	1:A:266:GLU:HG3	2.13	0.49
1:A:311:GLU:O	1:A:312:SER:C	2.49	0.49
1:A:466:LYS:HA	1:A:468:PRO:HD3	1.94	0.49
1:A:381:VAL:HG13	1:A:385:GLN:OE1	2.14	0.48
1:A:417:GLN:HG2	1:A:470:SER:HB2	1.94	0.48
1:A:135:LEU:O	1:A:139:LEU:HB2	2.14	0.48
1:A:156:PHE:CZ	1:A:285:GLU:HA	2.48	0.48
1:A:356:THR:O	1:A:360:CYS:N	2.36	0.48
1:A:341:TYR:CD2	1:A:341:TYR:N	2.82	0.48
1:A:197:ARG:HH11	1:A:458:ASN:HD21	1.62	0.48
1:A:425:GLU:HG3	1:A:429:ASN:OD1	2.13	0.48
1:A:444:LYS:C	1:A:447:PRO:HD2	2.34	0.48
1:A:512:ASP:O	1:A:516:LEU:HG	2.13	0.48
1:A:36:PHE:CG	1:A:137:LYS:HD3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:TYR:CZ	1:A:349:LEU:HD13	2.48	0.48
1:A:77:VAL:O	1:A:80:LEU:HB2	2.13	0.47
3:A:1006:11D:H26A	3:A:1006:11D:H23	1.55	0.47
1:A:218:ARG:NH1	1:A:451:ASP:OD1	2.47	0.47
1:A:32:GLN:NE2	1:A:110:PRO:HB3	2.29	0.47
1:A:329:MET:O	1:A:333:GLU:HG2	2.13	0.47
1:A:379:PRO:O	1:A:383:GLU:HB2	2.13	0.47
1:A:348:ARG:HG3	1:A:482:VAL:HB	1.95	0.47
1:A:505:GLU:CA	1:A:508:THR:HG22	2.40	0.47
1:A:14:LEU:O	1:A:18:ASN:ND2	2.47	0.47
1:A:221:GLN:O	1:A:224:PRO:HD3	2.15	0.47
1:A:199:LYS:HD2	1:A:199:LYS:HA	1.54	0.47
1:A:117:ARG:HH21	1:A:179:LEU:HD13	1.79	0.47
1:A:42:LEU:O	1:A:46:VAL:HG23	2.14	0.47
1:A:407:LEU:CD1	1:A:410:ARG:HH11	2.28	0.47
1:A:450:GLU:CG	1:A:451:ASP:N	2.78	0.47
1:A:416:PRO:HB2	1:A:497:TYR:CE2	2.50	0.47
1:A:25:ILE:O	1:A:29:GLN:HG3	2.14	0.47
1:A:77:VAL:HG22	1:A:80:LEU:HB2	1.96	0.47
1:A:203:LEU:HD12	1:A:203:LEU:O	2.15	0.47
1:A:480:SER:HG	1:A:482:VAL:HG22	1.80	0.46
1:A:480:SER:OG	1:A:482:VAL:CG2	2.63	0.46
1:A:525:LYS:HD3	2:A:1005:MYR:H31	1.95	0.46
1:A:439:LYS:HD3	1:A:439:LYS:HA	1.42	0.46
1:A:418:VAL:HA	1:A:469:VAL:CG2	2.44	0.46
1:A:570:GLU:HG3	1:A:571:GLU:H	1.80	0.46
1:A:161:TYR:O	1:A:165:PHE:HD2	1.98	0.46
1:A:399:GLY:O	1:A:403:PHE:HB2	2.16	0.46
1:A:397:GLN:O	1:A:398:LEU:HD12	2.15	0.46
1:A:525:LYS:HG2	1:A:551:PHE:CE2	2.51	0.46
1:A:42:LEU:HD22	1:A:73:LYS:CE	2.36	0.46
1:A:342:SER:OG	1:A:344:VAL:HG23	2.15	0.45
1:A:224:PRO:HB2	1:A:299:PRO:HD3	1.99	0.45
1:A:470:SER:O	1:A:473:VAL:HB	2.17	0.45
1:A:401:TYR:CE2	1:A:522:GLN:HG2	2.51	0.45
1:A:510:HIS:C	1:A:512:ASP:H	2.19	0.45
1:A:341:TYR:HD2	1:A:341:TYR:N	2.13	0.45
1:A:36:PHE:O	1:A:40:VAL:HG23	2.17	0.45
1:A:66:LEU:HD13	1:A:66:LEU:HA	1.59	0.45
1:A:545:LYS:O	1:A:549:ASP:HB2	2.16	0.45
1:A:120:VAL:CG1	1:A:174:LYS:HB3	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLN:O	1:A:197:ARG:C	2.55	0.45
1:A:391:ASN:HB2	2:A:1003:MYR:H101	1.98	0.45
1:A:224:PRO:HB2	1:A:298:MET:HA	1.98	0.45
1:A:149:PHE:CD1	1:A:149:PHE:C	2.91	0.45
1:A:251:LEU:HD23	2:A:1002:MYR:H91	1.98	0.45
1:A:405:ASN:ND2	1:A:405:ASN:N	2.64	0.45
1:A:217:ALA:O	1:A:221:GLN:HG3	2.17	0.45
1:A:238:LEU:HD22	3:A:1007:11D:H13B	1.99	0.44
1:A:67:HIS:HE1	1:A:249:ASP:OD2	2.01	0.44
1:A:480:SER:OG	1:A:482:VAL:HG22	2.18	0.44
1:A:488:PHE:HB3	2:A:1004:MYR:C9	2.33	0.44
1:A:530:VAL:O	1:A:533:VAL:N	2.50	0.44
1:A:348:ARG:HH12	2:A:1003:MYR:C1	2.31	0.44
1:A:325:VAL:O	1:A:329:MET:HG3	2.18	0.44
1:A:388:ILE:HD12	1:A:389:LYS:N	2.32	0.44
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.99	0.44
1:A:510:HIS:C	1:A:512:ASP:N	2.69	0.44
1:A:184:GLU:O	1:A:188:GLU:HG3	2.16	0.44
1:A:31:LEU:HD21	1:A:74:LEU:HD22	1.99	0.44
1:A:466:LYS:CA	1:A:466:LYS:HE3	2.35	0.44
1:A:90:CYS:O	1:A:98:ARG:HG3	2.18	0.44
2:A:1003:MYR:H132	2:A:1004:MYR:H51	1.98	0.44
1:A:156:PHE:HD2	1:A:156:PHE:N	2.11	0.44
1:A:458:ASN:OD1	1:A:484:ARG:NH1	2.49	0.44
1:A:570:GLU:C	1:A:572:GLY:N	2.69	0.44
1:A:544:LEU:C	1:A:546:ALA:N	2.66	0.44
2:A:1004:MYR:H101	2:A:1004:MYR:H71	1.55	0.43
1:A:178:LEU:O	1:A:179:LEU:C	2.56	0.43
1:A:195:LYS:HE3	1:A:195:LYS:HB2	1.71	0.43
1:A:282:PRO:O	1:A:283:LEU:C	2.56	0.43
1:A:540:THR:HB	1:A:542:GLU:CG	2.48	0.43
1:A:265:CYS:O	1:A:268:GLN:HG3	2.17	0.43
1:A:41:LYS:O	1:A:45:GLU:HG3	2.17	0.43
1:A:513:ILE:CG2	1:A:514:CYS:N	2.82	0.43
1:A:410:ARG:HG3	1:A:414:LYS:HE3	2.00	0.43
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.99	0.43
1:A:32:GLN:HE22	1:A:110:PRO:HB3	1.84	0.43
1:A:457:LEU:HD21	2:A:1004:MYR:H91	2.00	0.43
1:A:565:GLU:HA	1:A:565:GLU:OE2	2.18	0.43
1:A:26:ALA:O	1:A:30:TYR:HD2	2.02	0.43
1:A:404:GLN:NE2	1:A:428:ARG:HA	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1007:11D:O15	3:A:1007:11D:H10	2.19	0.43
1:A:149:PHE:HD1	1:A:149:PHE:C	2.22	0.43
1:A:156:PHE:HZ	1:A:284:LEU:HG	1.83	0.43
1:A:304:SER:OG	1:A:306:ALA:HB3	2.18	0.43
1:A:305:LEU:HB3	1:A:374:PHE:HZ	1.84	0.43
1:A:565:GLU:CA	1:A:565:GLU:OE2	2.66	0.43
1:A:234:LEU:O	1:A:238:LEU:HB2	2.18	0.43
3:A:1006:11D:H12A	3:A:1006:11D:H1	2.01	0.42
1:A:453:LEU:HD11	2:A:1004:MYR:H62	1.97	0.42
1:A:283:LEU:O	2:A:1002:MYR:H21	2.19	0.42
1:A:209:ARG:NH1	3:A:1006:11D:H26	2.34	0.42
1:A:340:ASP:O	1:A:447:PRO:HD3	2.19	0.42
1:A:269:ASP:OD2	1:A:269:ASP:N	2.53	0.42
1:A:296:ASP:OD1	1:A:297:GLU:N	2.52	0.42
1:A:332:TYR:OH	1:A:336:ARG:CZ	2.67	0.42
1:A:458:ASN:HA	1:A:484:ARG:NH1	2.34	0.42
1:A:152:PRO:O	1:A:156:PHE:CD2	2.73	0.42
1:A:163:ALA:O	1:A:167:GLU:HB2	2.19	0.42
1:A:117:ARG:HG2	1:A:123:MET:HE1	2.01	0.42
1:A:222:ARG:HG3	1:A:295:ASN:HD22	1.84	0.42
1:A:370:TYR:O	1:A:370:TYR:HD1	2.02	0.42
1:A:205:LYS:HE3	1:A:206:PHE:CE2	2.55	0.42
1:A:24:LEU:HD21	1:A:140:TYR:HB2	2.01	0.42
1:A:453:LEU:HD22	1:A:453:LEU:HA	1.80	0.42
1:A:446:MET:O	1:A:447:PRO:C	2.57	0.42
1:A:343:VAL:O	1:A:346:LEU:N	2.52	0.42
1:A:564:LYS:HG2	1:A:564:LYS:O	2.19	0.42
1:A:25:ILE:HG23	1:A:143:ALA:HB2	2.02	0.42
1:A:576:VAL:O	1:A:580:GLN:HB2	2.20	0.42
1:A:257:ARG:CZ	3:A:1007:11D:H26	2.51	0.41
1:A:259:ASP:HA	1:A:262:LYS:HE3	2.02	0.41
1:A:117:ARG:HA	1:A:118:PRO:HD3	1.89	0.41
1:A:384:PRO:HG3	1:A:485:ARG:HH22	1.85	0.41
1:A:486:PRO:O	1:A:487:CYS:C	2.58	0.41
1:A:551:PHE:O	1:A:555:VAL:HG22	2.20	0.41
1:A:74:LEU:O	1:A:77:VAL:HG12	2.18	0.41
1:A:178:LEU:HD22	1:A:182:LEU:HG	2.03	0.41
1:A:372:LYS:NZ	1:A:372:LYS:CB	2.84	0.41
1:A:220:SER:HB2	1:A:335:ALA:HB3	2.01	0.41
1:A:34:CYS:HA	1:A:35:PRO:HD3	1.90	0.41
1:A:360:CYS:O	1:A:362:ALA:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LYS:HA	1:A:73:LYS:HD2	1.67	0.41
1:A:127:PHE:O	1:A:131:GLU:HG2	2.19	0.41
1:A:161:TYR:O	1:A:165:PHE:CD2	2.73	0.41
1:A:295:ASN:OD1	1:A:339:PRO:HB2	2.21	0.41
1:A:330:PHE:HZ	1:A:377:PHE:CE1	2.38	0.41
1:A:430:LEU:HG	1:A:456:VAL:HG11	2.02	0.41
1:A:540:THR:C	1:A:542:GLU:H	2.24	0.41
1:A:572:GLY:O	1:A:573:LYS:C	2.59	0.41
1:A:10:ARG:O	1:A:14:LEU:HB2	2.20	0.41
1:A:441:PRO:O	1:A:442:GLU:C	2.59	0.41
1:A:179:LEU:O	1:A:180:PRO:C	2.55	0.41
1:A:191:ALA:O	1:A:195:LYS:N	2.42	0.41
1:A:429:ASN:HD22	1:A:432:LYS:HE2	1.86	0.41
1:A:328:GLY:HA2	3:A:1006:11D:H18	2.03	0.41
1:A:370:TYR:CD1	1:A:370:TYR:C	2.94	0.41
1:A:403:PHE:O	1:A:406:ALA:HB3	2.21	0.41
1:A:178:LEU:O	1:A:181:LYS:N	2.54	0.40
1:A:383:GLU:OE2	1:A:384:PRO:HG3	2.22	0.40
1:A:380:LEU:O	1:A:384:PRO:HD3	2.20	0.40
1:A:26:ALA:HB1	1:A:70:PHE:CE1	2.56	0.40
1:A:118:PRO:HD2	1:A:123:MET:SD	2.61	0.40
1:A:393:GLU:O	1:A:394:LEU:C	2.59	0.40
1:A:420:THR:N	1:A:421:PRO:CD	2.84	0.40
1:A:108:ASP:C	1:A:110:PRO:HD3	2.40	0.40
1:A:571:GLU:O	1:A:574:LYS:HG2	2.21	0.40
1:A:168:CYS:SG	1:A:178:LEU:N	2.94	0.40
1:A:266:GLU:C	1:A:267:ASN:ND2	2.75	0.40
1:A:458:ASN:HA	1:A:484:ARG:HH11	1.86	0.40
1:A:526:GLN:HB2	1:A:526:GLN:HE21	1.57	0.40
1:A:531:GLU:HA	1:A:534:LYS:HB2	2.04	0.40
1:A:450:GLU:HG3	1:A:451:ASP:N	2.34	0.40
1:A:428:ARG:NH2	1:A:526:GLN:HE22	2.12	0.40
1:A:81:ARG:O	1:A:85:GLY:HA2	2.22	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	574/585 (98%)	486 (85%)	78 (14%)	10 (2%)	11	21

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	ALA
1	A	467	THR
1	A	247	HIS
1	A	286	LYS
1	A	301	ASP
1	A	361	CYS
1	A	442	GLU
1	A	447	PRO
1	A	558	CYS
1	A	366	PRO

### 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	506/511 (99%)	430 (85%)	76 (15%)	3	6

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS

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Mol	Chain	Res	Type
1	A	5	SER
1	A	7	VAL
1	A	13	ASP
1	A	20	LYS
1	A	41	LYS
1	A	60	GLU
1	A	66	LEU
1	A	77	VAL
1	A	82	GLU
1	A	87	MET
1	A	94	GLN
1	A	108	ASP
1	A	114	ARG
1	A	116	VAL
1	A	130	ASN
1	A	149	PHE
1	A	154	LEU
1	A	179	LEU
1	A	185	LEU
1	A	197	ARG
1	A	199	LYS
1	A	218	ARG
1	A	220	SER
1	A	222	ARG
1	A	235	VAL
1	A	241	VAL
1	A	255	ASP
1	A	269	ASP
1	A	273	SER
1	A	286	LYS
1	A	294	GLU
1	A	310	VAL
1	A	324	ASP
1	A	334	TYR
1	A	344	VAL
1	A	345	LEU
1	A	356	THR
1	A	357	LEU
1	A	367	HIS
1	A	372	LYS
1	A	373	VAL
1	A	394	LEU

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Mol	Chain	Res	Type
1	A	405	ASN
1	A	407	LEU
1	A	417	GLN
1	A	427	SER
1	A	433	VAL
1	A	450	GLU
1	A	451	ASP
1	A	453	LEU
1	A	455	VAL
1	A	463	LEU
1	A	466	LYS
1	A	469	VAL
1	A	471	ASP
1	A	474	THR
1	A	480	SER
1	A	484	ARG
1	A	485	ARG
1	A	491	LEU
1	A	508	THR
1	A	515	THR
1	A	527	THR
1	A	538	LYS
1	A	541	LYS
1	A	542	GLU
1	A	554	PHE
1	A	556	GLU
1	A	559	CYS
1	A	564	LYS
1	A	567	CYS
1	A	570	GLU
1	A	571	GLU
1	A	574	LYS
1	A	580	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	67	HIS
1	A	105	HIS
1	A	130	ASN
1	A	146	HIS

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Mol	Chain	Res	Type
1	A	242	HIS
1	A	267	ASN
1	A	268	GLN
1	A	295	ASN
1	A	338	HIS
1	A	390	GLN
1	A	391	ASN
1	A	397	GLN
1	A	405	ASN
1	A	429	ASN
1	A	440	HIS
1	A	459	GLN
1	A	483	ASN
1	A	522	GLN
1	A	526	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	MYR	A	1002	-	12,15,15	0.18	0	11,15,15	0.91	0
2	MYR	A	1003	-	12,15,15	0.32	0	11,15,15	0.73	0
2	MYR	A	1004	-	11,11,15	0.79	0	10,10,15	2.78	7 (70%)
2	MYR	A	1005	-	12,15,15	0.24	0	11,15,15	0.65	0
3	11D	A	1006	-	28,31,31	2.95	3 (10%)	37,41,41	1.83	9 (24%)
3	11D	A	1007	-	28,31,31	3.69	2 (7%)	37,41,41	1.65	6 (16%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MYR	A	1002	-	-	0/11/13/13	0/0/0/0
2	MYR	A	1003	-	-	0/11/13/13	0/0/0/0
2	MYR	A	1004	-	-	0/9/9/13	0/0/0/0
2	MYR	A	1005	-	-	0/11/13/13	0/0/0/0
3	11D	A	1006	-	-	0/23/25/25	0/2/2/2
3	11D	A	1007	-	-	0/23/25/25	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1007	11D	C4-S14	-18.63	1.58	1.78
3	A	1006	11D	C4-S14	-14.66	1.62	1.78
3	A	1007	11D	C4-C3	-3.88	1.37	1.43
3	A	1006	11D	S14-N17	-2.62	1.58	1.61
3	A	1006	11D	C5-C4	2.84	1.40	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1007	11D	O15-S14-O16	-6.61	111.09	119.55
3	A	1006	11D	O15-S14-O16	-6.54	111.18	119.55
2	A	1004	MYR	C8-C7-C6	-5.03	88.56	114.45
2	A	1004	MYR	C11-C10-C9	-3.67	95.54	114.45
3	A	1006	11D	C5-C4-C3	-2.88	118.14	120.98
2	A	1004	MYR	C9-C8-C7	-2.85	99.79	114.45
2	A	1004	MYR	C7-C6-C5	-2.82	99.92	114.45
2	A	1004	MYR	C13-C12-C11	-2.77	88.30	115.40
2	A	1004	MYR	C12-C11-C10	-2.65	100.78	114.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1007	11D	C10-C3-C4	-2.57	120.33	123.63
3	A	1007	11D	C3-C4-S14	-2.55	118.42	121.56
3	A	1006	11D	C10-C3-C4	-2.39	120.57	123.63
2	A	1004	MYR	C10-C9-C8	-2.14	103.43	114.45
3	A	1007	11D	C10-C3-C2	2.10	121.15	118.49
3	A	1006	11D	C13-N11-C7	2.13	120.34	114.01
3	A	1007	11D	C5-C4-S14	2.21	119.75	116.98
3	A	1006	11D	C3-C4-S14	2.22	124.29	121.56
3	A	1007	11D	C13-N11-C7	2.50	121.42	114.01
3	A	1006	11D	C2-C7-N11	2.66	121.97	118.17
3	A	1006	11D	C4-C3-C2	2.73	120.92	117.91
3	A	1006	11D	C12-N11-C7	2.98	122.85	114.01
3	A	1006	11D	C4-S14-N17	3.06	112.78	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002	MYR	8	0
2	A	1003	MYR	16	0
2	A	1004	MYR	19	0
2	A	1005	MYR	3	0
3	A	1006	11D	8	0
3	A	1007	11D	8	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	578/585 (98%)	0.27	14 (2%) 59 52	36, 60, 90, 110	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	559	CYS	3.6
1	A	553	ALA	3.6
1	A	557	LYS	2.7
1	A	138	TYR	2.5
1	A	442	GLU	2.5
1	A	582	ALA	2.3
1	A	197	ARG	2.3
1	A	495	GLU	2.2
1	A	580	GLN	2.2
1	A	494	ASP	2.1
1	A	502	PHE	2.0
1	A	294	GLU	2.0
1	A	204	GLN	2.0
1	A	498	VAL	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 carbohydrates 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. 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, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MYR	A	1003	16/16	0.87	0.33	4.78	53,63,69,70	0
2	MYR	A	1005	16/16	0.89	0.41	3.17	61,67,82,82	0
2	MYR	A	1004	12/16	0.89	0.25	2.10	54,55,72,75	0
3	11D	A	1006	30/30	0.91	0.26	1.89	54,66,74,80	0
3	11D	A	1007	30/30	0.92	0.24	1.30	52,65,78,83	0
2	MYR	A	1002	16/16	0.88	0.23	1.03	52,55,62,62	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.