



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:34 PM UTC

PDB ID : 3R5A / pdb_00003r5a
Title : Pseudomonas aeruginosa DapD (PA3666) in complex with D-2-aminopimelate
Authors : Sandalova, T.; Schnell, R.; Schneider, G.
Deposited on : 2011-03-18
Resolution : 1.89 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

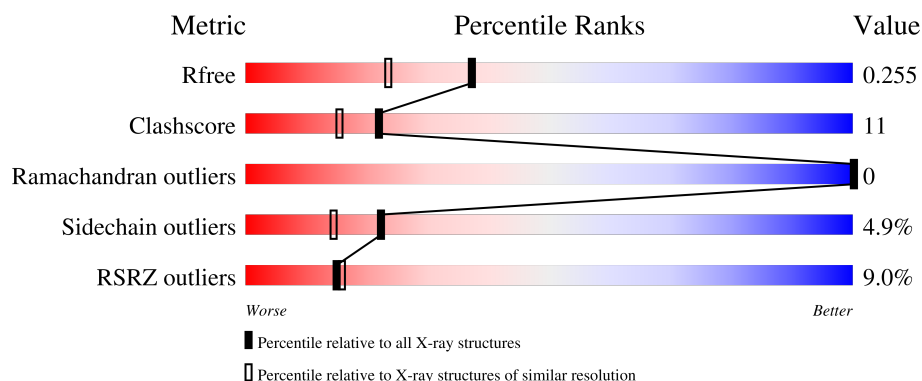
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	347	8% 80% 12% 7%
1	B	347	10% 73% 16% 8%
1	C	347	8% 69% 22% 7%
1	D	347	8% 78% 12% 7%
1	E	347	8% 73% 14% 10%

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Mol	Chain	Length	Quality of chain
1	F	347	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	E	345	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15195 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 Tetrahydrodipicolinate N-succinyletransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2364	1506	406	445	7			
1	B	318	Total	C	N	O	S	0	0	0
			2342	1492	403	440	7			
1	C	321	Total	C	N	O	S	0	0	0
			2367	1506	407	447	7			
1	D	321	Total	C	N	O	S	0	0	0
			2364	1506	406	445	7			
1	E	311	Total	C	N	O	S	0	0	0
			2288	1464	387	430	7			
1	F	321	Total	C	N	O	S	0	0	0
			2367	1506	407	447	7			

There are 18 discrepancies between the modelled and reference sequences:

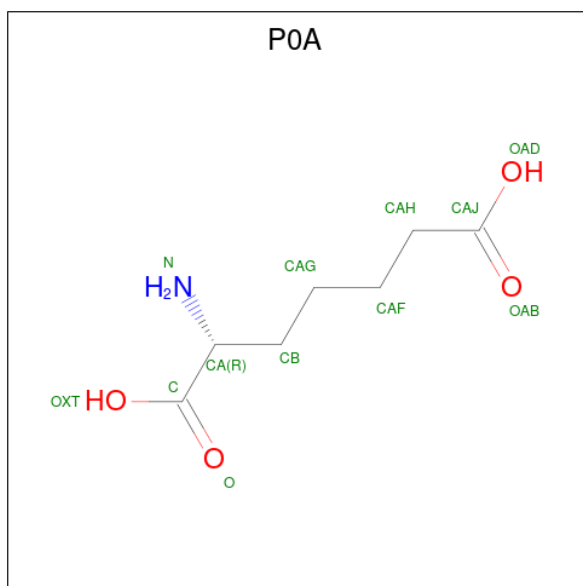
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9Z9H2
A	-1	SER	-	expression tag	UNP Q9Z9H2
A	0	HIS	-	expression tag	UNP Q9Z9H2
B	-2	GLY	-	expression tag	UNP Q9Z9H2
B	-1	SER	-	expression tag	UNP Q9Z9H2
B	0	HIS	-	expression tag	UNP Q9Z9H2
C	-2	GLY	-	expression tag	UNP Q9Z9H2
C	-1	SER	-	expression tag	UNP Q9Z9H2
C	0	HIS	-	expression tag	UNP Q9Z9H2
D	-2	GLY	-	expression tag	UNP Q9Z9H2
D	-1	SER	-	expression tag	UNP Q9Z9H2
D	0	HIS	-	expression tag	UNP Q9Z9H2
E	-2	GLY	-	expression tag	UNP Q9Z9H2
E	-1	SER	-	expression tag	UNP Q9Z9H2
E	0	HIS	-	expression tag	UNP Q9Z9H2
F	-2	GLY	-	expression tag	UNP Q9Z9H2
F	-1	SER	-	expression tag	UNP Q9Z9H2

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	-	expression tag	UNP Q9Z9H2

- Molecule 2 is (2R)-2-aminoheptanedioic acid (CCD ID: P0A) (formula: $C_7H_{13}NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			12	7	1	4		
2	A	1	Total	C	N	O	0	0
			12	7	1	4		
2	C	1	Total	C	N	O	0	0
			12	7	1	4		
2	D	1	Total	C	N	O	0	0
			12	7	1	4		
2	E	1	Total	C	N	O	0	0
			12	7	1	4		
2	F	1	Total	C	N	O	0	0
			12	7	1	4		

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



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

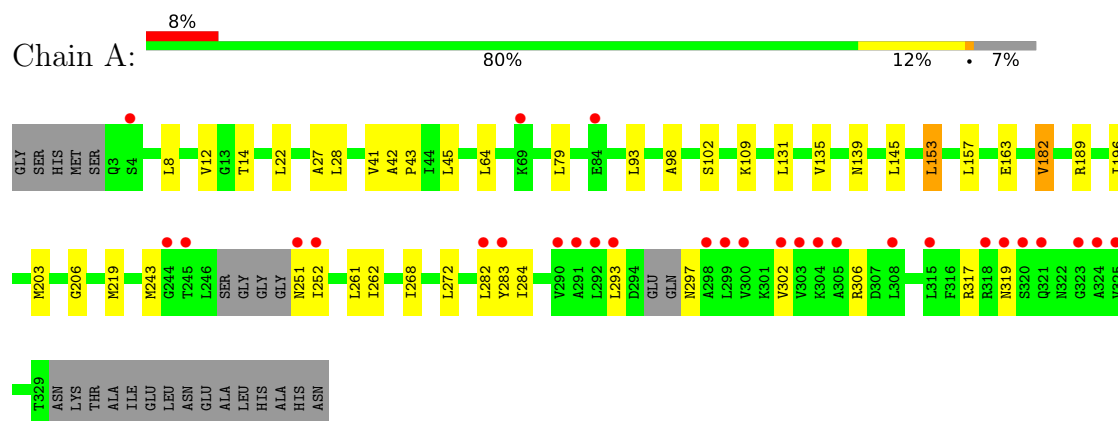
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	175	Total	O	0	0
			175	175		
4	B	192	Total	O	0	0
			192	192		
4	C	143	Total	O	0	0
			143	143		
4	D	172	Total	O	0	0
			172	172		
4	E	212	Total	O	0	0
			212	212		
4	F	125	Total	O	0	0
			125	125		

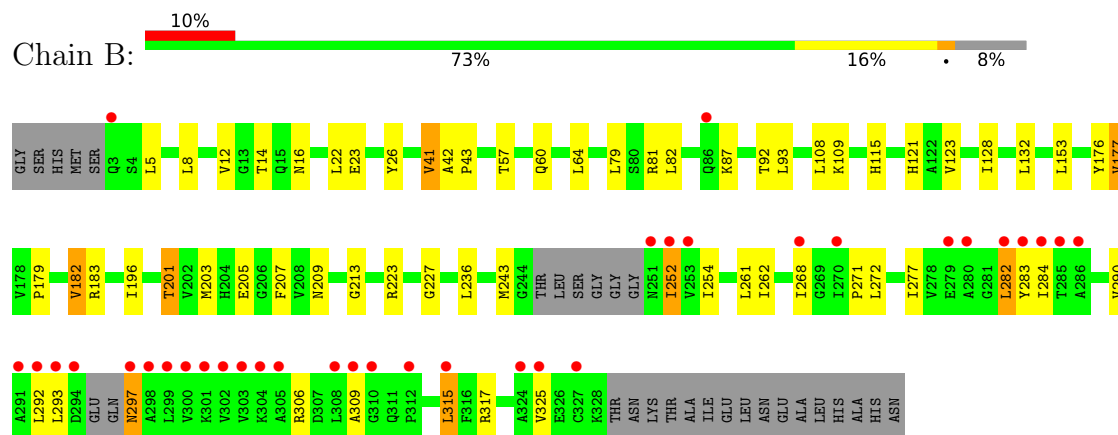
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

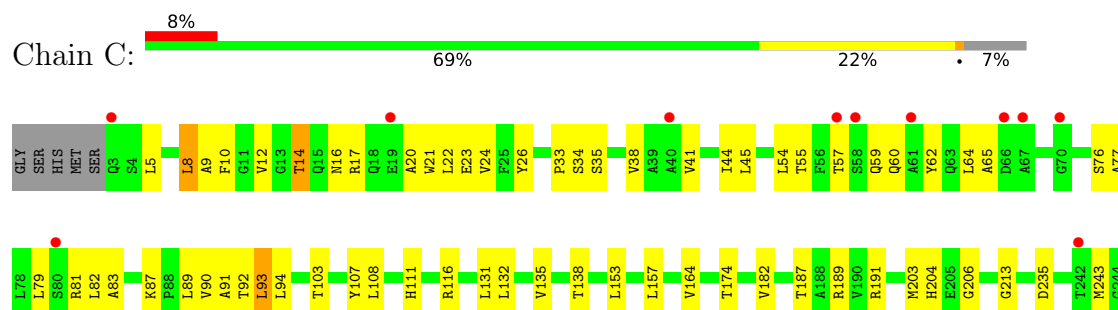
- Molecule 1: Tetrahydrodipicolinate N-succinyletransferase



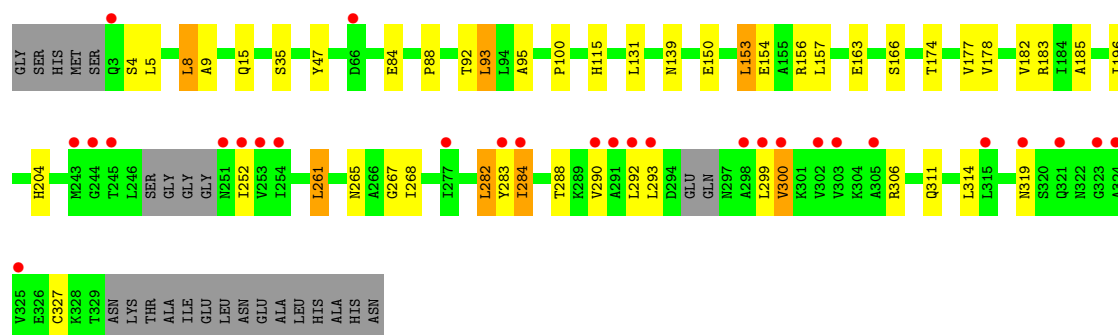
- Molecule 1: Tetrahydrodipicolinate N-succinyletransferase



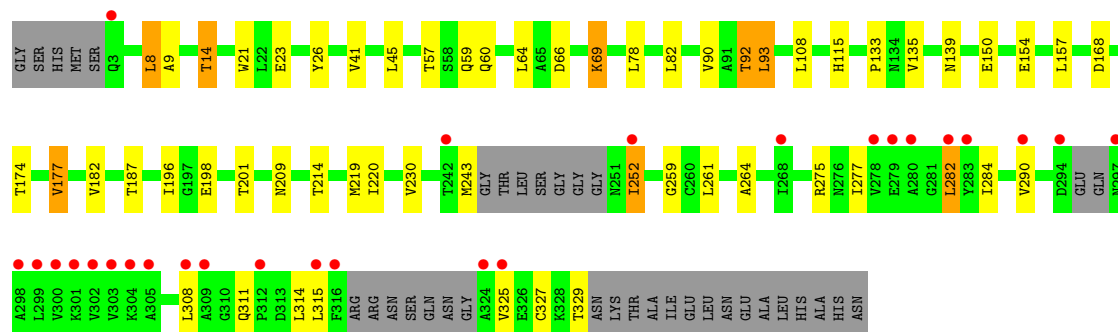
- Molecule 1: Tetrahydrodipicolinate N-succinyletransferase



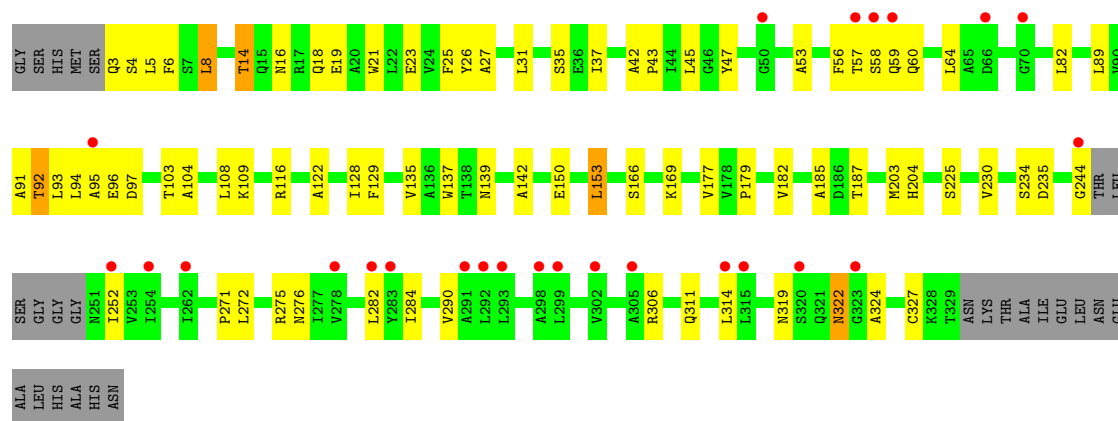
Chain D: 8% 78% 12% 2%



Chain E:



Chain F: 7% 69% 22% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.34Å 101.93Å 135.69Å 90.00° 89.97° 90.00°	Depositor
Resolution (Å)	21.83 – 1.89 21.83 – 1.89	Depositor EDS
% Data completeness (in resolution range)	96.1 (21.83-1.89) 98.5 (21.83-1.89)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.196 , 0.235 (Not available) , 0.255	Depositor DCC
R_{free} test set	9008 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.277 for h,-k,-l	Xtriage
Reported twinning fraction	0.509 for H, K, L 0.491 for h,-k,-l	Depositor
Outliers	0 of 178640 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15195	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: GOL, POA

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	0.55	0/2400	0.74	0/3257
1	B	0.56	0/2378	0.76	1/3226 (0.0%)
1	C	0.55	0/2404	0.76	1/3263 (0.0%)
1	D	0.54	0/2400	0.77	1/3257 (0.0%)
1	E	0.57	0/2323	0.76	0/3153
1	F	0.57	0/2404	0.78	1/3263 (0.0%)
All	All	0.56	0/14309	0.76	4/19419 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	213	GLY	N-CA-C	6.99	119.42	110.20
1	B	213	GLY	N-CA-C	5.46	117.27	110.29
1	D	177	VAL	N-CA-C	5.30	114.85	108.06
1	F	177	VAL	N-CA-C	5.29	115.34	108.35

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	2364	0	2420	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2342	0	2395	53	0
1	C	2367	0	2417	73	0
1	D	2364	0	2420	33	0
1	E	2288	0	2344	54	0
1	F	2367	0	2417	69	0
2	A	24	0	22	4	0
2	C	12	0	11	2	0
2	D	12	0	11	1	0
2	E	12	0	11	2	0
2	F	12	0	11	0	0
3	A	6	0	8	2	0
3	E	6	0	8	5	0
4	A	175	0	0	1	0
4	B	192	0	0	4	0
4	C	143	0	0	3	0
4	D	172	0	0	2	0
4	E	212	0	0	8	0
4	F	125	0	0	4	0
All	All	15195	0	14495	306	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 11.

All (306) 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:F:45:LEU:HD21	1:F:64:LEU:HD11	1.34	1.06
1:E:201:THR:HB	4:E:996:HOH:O	1.63	0.98
1:C:14:THR:HG22	1:C:87:LYS:NZ	1.79	0.98
1:B:14:THR:HG22	1:B:87:LYS:NZ	1.80	0.96
1:F:57:THR:HG22	1:F:60:GLN:NE2	1.81	0.95
1:B:201:THR:HG22	4:B:713:HOH:O	1.67	0.93
1:C:14:THR:HG23	1:C:23:GLU:OE1	1.70	0.91
1:B:41:VAL:HG22	1:B:64:LEU:HD22	1.54	0.90
1:B:268:ILE:HG21	1:B:272:LEU:HD11	1.55	0.89
3:E:345:GOL:H12	1:F:139:ASN:HB3	1.56	0.88
1:F:14:THR:HG22	1:F:23:GLU:OE1	1.73	0.88
1:F:252:ILE:HD12	1:F:252:ILE:O	1.74	0.87
1:E:41:VAL:HG13	1:E:64:LEU:HD22	1.56	0.87
1:C:22:LEU:HD22	1:C:135:VAL:HG22	1.59	0.85
3:E:345:GOL:H12	1:F:139:ASN:CB	2.08	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:THR:HG22	1:C:87:LYS:HZ1	1.44	0.82
1:D:284:ILE:HG22	1:D:288:THR:HG21	1.61	0.82
1:B:14:THR:HG23	1:B:23:GLU:OE2	1.79	0.82
1:E:311:GLN:HB2	1:E:314:LEU:HD11	1.63	0.81
1:C:116:ARG:HD3	4:C:910:HOH:O	1.82	0.79
1:F:94:LEU:HD11	1:F:108:LEU:HD23	1.65	0.78
1:A:252:ILE:HD12	1:A:252:ILE:O	1.84	0.78
1:B:177:VAL:HB	4:B:988:HOH:O	1.83	0.78
1:C:8:LEU:HB2	1:C:93:LEU:HD22	1.65	0.77
1:D:311:GLN:HB2	1:D:314:LEU:HD11	1.66	0.77
1:E:41:VAL:HG13	1:E:64:LEU:CD2	2.15	0.76
1:B:14:THR:HG22	1:B:87:LYS:HZ1	1.49	0.76
1:E:14:THR:HG22	1:E:23:GLU:OE2	1.87	0.74
1:C:14:THR:HG22	1:C:87:LYS:HZ3	1.51	0.74
1:F:5:LEU:HD22	1:F:95:ALA:HA	1.69	0.74
1:F:137:TRP:CZ3	1:F:142:ALA:HB2	2.21	0.74
1:D:8:LEU:HD12	1:D:9:ALA:N	2.05	0.72
1:F:97:ASP:OD1	1:F:109:LYS:NZ	2.21	0.71
1:B:293:LEU:HD13	1:B:297:ASN:HB2	1.72	0.71
1:B:14:THR:HG22	1:B:87:LYS:HZ3	1.52	0.71
1:C:189:ARG:HD3	1:C:206:GLY:O	1.91	0.71
1:F:26:TYR:OH	1:F:92:THR:HG21	1.91	0.71
1:E:264:ALA:HB2	4:E:451:HOH:O	1.91	0.70
1:C:41:VAL:HG22	1:C:64:LEU:HD22	1.72	0.70
1:B:179:PRO:HG2	1:B:182:VAL:HG21	1.74	0.69
1:F:42:ALA:HB3	1:F:43:PRO:HD3	1.74	0.69
1:F:45:LEU:CD2	1:F:64:LEU:HD11	2.17	0.67
1:A:41:VAL:HG22	1:A:64:LEU:HD22	1.77	0.67
1:B:26:TYR:OH	1:B:92:THR:HG21	1.93	0.67
1:E:21:TRP:CH2	1:E:90:VAL:HG13	2.30	0.67
1:A:41:VAL:HG12	1:A:45:LEU:HD12	1.77	0.67
1:E:209:ASN:HD21	2:E:351:P0A:HAF	1.58	0.66
1:E:201:THR:HG21	1:E:219:MET:CE	2.24	0.66
1:E:8:LEU:HB2	1:E:93:LEU:HD22	1.77	0.66
3:E:345:GOL:C1	1:F:139:ASN:HB3	2.24	0.66
1:F:47:TYR:OH	1:F:95:ALA:HB2	1.97	0.64
1:D:182:VAL:HG11	1:D:196:ILE:HG21	1.78	0.64
1:F:92:THR:OG1	1:F:108:LEU:HD22	1.98	0.64
1:B:290:VAL:HG13	1:B:325:VAL:HG12	1.79	0.64
1:C:294:ASP:O	1:C:296:GLN:O	2.16	0.64
1:F:57:THR:CG2	1:F:60:GLN:NE2	2.58	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:LEU:CD2	1:B:284:ILE:HG23	2.29	0.62
1:A:297:ASN:N	4:A:1009:HOH:O	2.33	0.62
1:A:139:ASN:HB2	3:A:346:GOL:H12	1.80	0.62
1:D:100:PRO:O	1:D:131:LEU:HD11	1.99	0.62
1:C:5:LEU:HD22	1:C:94:LEU:O	1.99	0.62
1:F:94:LEU:CD1	1:F:108:LEU:HD23	2.29	0.62
1:F:311:GLN:HB2	1:F:314:LEU:HD11	1.82	0.61
1:E:57:THR:H	1:E:60:GLN:HE21	1.49	0.61
1:B:121:HIS:NE2	1:B:227:GLY:O	2.31	0.60
1:A:182:VAL:HG22	1:A:196:ILE:HG22	1.83	0.60
1:B:205:GLU:OE2	1:C:189:ARG:NH2	2.34	0.60
1:F:45:LEU:HD21	1:F:64:LEU:CD1	2.23	0.60
1:F:57:THR:CG2	1:F:60:GLN:HE21	2.15	0.60
1:F:275:ARG:NH1	4:F:362:HOH:O	2.35	0.60
1:C:92:THR:C	1:C:93:LEU:HD23	2.27	0.59
1:F:14:THR:CG2	1:F:23:GLU:OE1	2.49	0.59
1:A:182:VAL:CG2	1:A:196:ILE:HG22	2.32	0.59
1:E:282:LEU:HD11	1:E:325:VAL:HG13	1.85	0.59
1:E:115:HIS:CD2	1:E:177:VAL:HG11	2.37	0.59
1:E:282:LEU:HD22	1:E:284:ILE:HG23	1.83	0.59
1:C:57:THR:HG23	1:C:60:GLN:NE2	2.18	0.59
1:B:252:ILE:HD13	1:B:254:ILE:HD11	1.85	0.58
1:C:90:VAL:HG22	1:C:92:THR:HG23	1.86	0.58
1:C:103:THR:OG1	1:C:131:LEU:O	2.20	0.58
1:A:14:THR:HG22	1:A:22:LEU:HD12	1.85	0.58
1:B:64:LEU:HB3	1:B:79:LEU:HD13	1.85	0.58
2:D:351:P0A:HAG	1:F:203:MET:HE2	1.85	0.57
1:F:235:ASP:HB3	4:F:377:HOH:O	2.04	0.57
1:A:268:ILE:HG21	1:A:272:LEU:HD11	1.85	0.57
1:F:252:ILE:HD13	1:F:271:PRO:HD3	1.87	0.57
1:B:128:ILE:HG12	1:B:132:LEU:HD11	1.86	0.57
1:B:207:PHE:HB3	1:B:223:ARG:HG2	1.85	0.57
1:C:108:LEU:O	1:C:108:LEU:HD12	2.05	0.57
1:F:16:ASN:OD1	1:F:18:GLN:N	2.38	0.57
1:E:115:HIS:NE2	1:E:177:VAL:HG11	2.19	0.57
1:E:139:ASN:HB3	3:E:345:GOL:H31	1.87	0.56
1:F:135:VAL:HG11	1:F:142:ALA:HB1	1.87	0.56
1:B:92:THR:C	1:B:93:LEU:HD23	2.31	0.56
1:B:277:ILE:HG21	1:B:315:LEU:HD12	1.86	0.56
1:D:292:LEU:O	1:D:300:VAL:HG22	2.06	0.56
1:B:115:HIS:CE1	1:B:177:VAL:HG11	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:16:ASN:HD21	1:C:20:ALA:HB3	1.71	0.55
1:A:293:LEU:HD13	1:A:297:ASN:HB2	1.88	0.54
1:A:41:VAL:HG12	1:A:45:LEU:CD1	2.36	0.54
1:C:10:PHE:CE1	1:C:79:LEU:HD11	2.43	0.54
1:F:92:THR:HG21	1:F:104:ALA:HB1	1.89	0.54
1:C:16:ASN:ND2	1:C:20:ALA:HB3	2.22	0.54
1:D:182:VAL:HG11	1:D:196:ILE:CG2	2.37	0.54
1:B:201:THR:CG2	4:B:713:HOH:O	2.40	0.54
1:C:44:ILE:CD1	1:C:64:LEU:HD23	2.37	0.54
1:C:189:ARG:HD2	1:C:204:HIS:O	2.07	0.54
1:E:115:HIS:NE2	1:E:177:VAL:CG1	2.71	0.54
1:C:14:THR:HG21	4:C:582:HOH:O	2.07	0.54
1:C:14:THR:CG2	1:C:87:LYS:HZ1	2.16	0.54
1:E:26:TYR:OH	1:E:92:THR:HG21	2.06	0.54
1:E:14:THR:CG2	1:E:23:GLU:OE2	2.55	0.53
1:C:8:LEU:HD23	1:C:33:PRO:HG3	1.90	0.53
1:C:45:LEU:HB3	1:C:54:LEU:HD13	1.90	0.53
1:F:45:LEU:HG	1:F:64:LEU:HD21	1.91	0.53
1:B:92:THR:OG1	1:B:108:LEU:HD22	2.09	0.53
1:C:41:VAL:HG13	1:C:64:LEU:CD2	2.39	0.53
1:F:47:TYR:HB2	1:F:93:LEU:HD12	1.90	0.53
1:F:16:ASN:OD1	1:F:16:ASN:C	2.51	0.53
1:B:236:LEU:HG	1:B:262:ILE:HD12	1.91	0.53
1:C:24:VAL:HG11	1:C:90:VAL:CG1	2.39	0.53
1:E:201:THR:HG23	1:E:219:MET:HA	1.90	0.53
1:F:47:TYR:CB	1:F:93:LEU:HD12	2.38	0.52
1:A:41:VAL:HG13	1:A:64:LEU:CD2	2.39	0.52
1:D:84:GLU:HB2	4:D:717:HOH:O	2.08	0.52
1:A:302:VAL:HG11	1:C:321:GLN:HE21	1.73	0.52
1:E:8:LEU:HD12	1:E:8:LEU:C	2.34	0.52
1:F:225:SER:HB3	1:F:244:GLY:HA2	1.90	0.52
1:C:14:THR:CG2	1:C:23:GLU:OE1	2.53	0.52
1:C:62:TYR:CZ	1:C:83:ALA:HB1	2.45	0.52
1:E:59:GLN:NE2	4:E:949:HOH:O	2.43	0.52
1:E:290:VAL:HG11	1:E:327:CYS:HB2	1.91	0.52
1:C:21:TRP:HZ2	1:C:55:THR:HG22	1.75	0.52
1:D:8:LEU:HD13	1:D:93:LEU:CD2	2.40	0.51
1:C:45:LEU:CD2	1:C:91:ALA:HB3	2.40	0.51
1:C:22:LEU:HD22	1:C:135:VAL:CG2	2.35	0.51
1:D:15:GLN:HE21	1:D:88:PRO:HG3	1.75	0.51
1:D:174:THR:HG22	1:D:178:VAL:HG22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:8:LEU:C	1:F:8:LEU:HD12	2.35	0.51
1:D:267:GLY:HA3	1:D:283:TYR:CE2	2.46	0.51
1:D:267:GLY:HA3	1:D:283:TYR:CD2	2.46	0.50
1:A:282:LEU:HD23	1:A:283:TYR:N	2.27	0.50
1:B:57:THR:H	1:B:60:GLN:HE21	1.59	0.50
1:E:21:TRP:CZ2	1:E:90:VAL:HG13	2.46	0.50
1:E:201:THR:HG21	1:E:219:MET:HE2	1.92	0.50
1:C:277:ILE:HB	1:C:315:LEU:HD12	1.93	0.50
1:C:64:LEU:HB3	1:C:79:LEU:HD13	1.93	0.50
1:A:41:VAL:CG1	1:A:45:LEU:CD1	2.90	0.49
1:A:139:ASN:ND2	1:A:163:GLU:H	2.10	0.49
1:B:271:PRO:HD2	1:B:309:ALA:HB1	1.94	0.49
1:B:292:LEU:HD12	1:B:293:LEU:H	1.77	0.49
1:C:44:ILE:HD12	1:C:64:LEU:HD23	1.94	0.49
1:B:177:VAL:HG12	4:B:371:HOH:O	2.13	0.49
1:B:182:VAL:HG22	1:B:196:ILE:HG22	1.95	0.49
1:E:259:GLY:O	1:E:275:ARG:HA	2.13	0.49
1:C:45:LEU:HD21	1:C:91:ALA:CB	2.43	0.49
1:F:57:THR:OG1	1:F:58:SER:N	2.46	0.48
1:C:191:ARG:HH12	2:C:351:P0A:HAH	1.78	0.48
1:B:282:LEU:C	1:B:282:LEU:HD23	2.38	0.48
1:C:138:THR:HG22	1:C:164:VAL:HG22	1.96	0.48
1:D:154:GLU:HA	1:D:157:LEU:HD12	1.95	0.48
1:E:78:LEU:O	1:E:82:LEU:HD22	2.13	0.48
1:E:201:THR:HG22	4:E:587:HOH:O	2.13	0.48
1:C:82:LEU:HD22	1:C:89:LEU:HD22	1.95	0.48
1:D:290:VAL:HG11	1:D:327:CYS:HB2	1.96	0.48
1:F:16:ASN:OD1	1:F:19:GLU:N	2.46	0.48
1:C:45:LEU:HD21	1:C:91:ALA:HB3	1.95	0.48
1:C:252:ILE:O	1:C:252:ILE:HD12	2.13	0.48
1:A:203:MET:SD	2:A:345:P0A:HAHA	2.54	0.48
1:A:219:MET:HE1	2:A:345:P0A:HAH	1.95	0.48
1:A:282:LEU:HD22	1:A:284:ILE:HG23	1.95	0.48
1:B:282:LEU:HD12	1:B:317:ARG:HA	1.96	0.48
1:D:261:LEU:C	1:D:261:LEU:HD13	2.39	0.48
1:E:14:THR:HG23	1:E:23:GLU:HG3	1.95	0.48
1:E:214:THR:HG22	1:E:230:VAL:HB	1.95	0.48
1:B:203:MET:HE1	1:C:191:ARG:NH2	2.29	0.47
1:B:282:LEU:HD21	1:B:284:ILE:HG23	1.94	0.47
1:C:77:ALA:O	1:C:81:ARG:HG2	2.13	0.47
1:F:57:THR:HG23	1:F:60:GLN:H	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:VAL:HG13	1:A:64:LEU:HD21	1.96	0.47
1:F:322:ASN:OD1	1:F:324:ALA:HB2	2.13	0.47
1:C:107:TYR:CE1	1:C:132:LEU:HD22	2.48	0.47
1:E:182:VAL:HG11	1:E:196:ILE:HG21	1.96	0.47
1:B:282:LEU:CD2	1:B:282:LEU:C	2.87	0.47
1:D:139:ASN:ND2	1:D:163:GLU:H	2.12	0.47
1:C:34:SER:O	1:C:38:VAL:HG23	2.15	0.47
1:C:57:THR:HG23	1:C:60:GLN:HE21	1.78	0.47
1:C:174:THR:HG21	1:C:187:THR:CG2	2.44	0.47
1:E:59:GLN:NE2	4:E:601:HOH:O	2.47	0.47
1:E:308:LEU:HD13	1:E:327:CYS:SG	2.55	0.47
1:F:93:LEU:C	1:F:94:LEU:HD23	2.40	0.47
1:F:8:LEU:HD21	1:F:37:ILE:HG22	1.97	0.47
1:E:252:ILE:HD12	1:E:252:ILE:O	2.15	0.47
1:B:41:VAL:HG22	1:B:64:LEU:CD2	2.35	0.46
1:B:203:MET:HE2	2:C:351:P0A:HAFA	1.98	0.46
1:F:290:VAL:HG11	1:F:327:CYS:HB2	1.97	0.46
1:E:26:TYR:OH	1:E:92:THR:CG2	2.64	0.46
1:E:78:LEU:O	1:E:82:LEU:CD2	2.64	0.46
1:E:174:THR:OG1	1:E:187:THR:HB	2.15	0.46
1:B:203:MET:HE1	1:C:191:ARG:CZ	2.45	0.46
1:A:261:LEU:HD23	1:A:262:ILE:N	2.31	0.46
1:B:8:LEU:C	1:B:8:LEU:HD12	2.40	0.46
1:B:115:HIS:CE1	1:B:177:VAL:CG1	2.99	0.46
1:F:272:LEU:HD22	1:F:276:ASN:HD22	1.80	0.46
1:A:302:VAL:HG11	1:C:321:GLN:NE2	2.30	0.46
1:B:109:LYS:HE2	1:B:123:VAL:CG2	2.46	0.46
1:B:261:LEU:HD23	1:B:262:ILE:N	2.31	0.46
1:C:82:LEU:HD22	1:C:89:LEU:CD2	2.46	0.46
1:D:5:LEU:HD22	1:D:95:ALA:HA	1.98	0.46
1:F:21:TRP:O	1:F:103:THR:HG21	2.16	0.46
1:B:128:ILE:HG12	1:B:132:LEU:CD1	2.46	0.45
2:E:351:P0A:HAG	2:E:351:P0A:OAB	2.16	0.45
1:F:8:LEU:HD21	1:F:37:ILE:CG2	2.47	0.45
1:C:243:MET:HE3	1:C:269:GLY:CA	2.45	0.45
1:E:57:THR:H	1:E:60:GLN:NE2	2.12	0.45
1:F:93:LEU:O	1:F:94:LEU:HD23	2.16	0.45
1:A:64:LEU:HB3	1:A:79:LEU:HD13	1.98	0.45
1:B:282:LEU:HD23	1:B:283:TYR:N	2.31	0.45
1:C:41:VAL:HG13	1:C:64:LEU:HD21	1.98	0.45
1:F:185:ALA:HB1	1:F:204:HIS:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:SER:HA	1:F:187:THR:HG23	1.99	0.45
1:A:98:ALA:O	1:A:109:LYS:NZ	2.44	0.45
1:B:176:TYR:C	1:B:177:VAL:CG1	2.90	0.45
1:C:243:MET:HE3	1:C:269:GLY:HA3	1.99	0.45
1:D:92:THR:C	1:D:93:LEU:HD23	2.41	0.45
1:D:293:LEU:HD22	1:D:299:LEU:HA	1.97	0.45
1:F:6:PHE:O	1:F:31:LEU:HD12	2.17	0.45
1:B:57:THR:H	1:B:60:GLN:NE2	2.15	0.44
1:C:65:ALA:O	1:C:76:SER:OG	2.35	0.44
1:F:14:THR:HG23	1:F:23:GLU:HG3	1.99	0.44
1:E:41:VAL:HG12	1:E:45:LEU:HD12	2.00	0.44
1:C:282:LEU:HD22	1:C:284:ILE:CG2	2.48	0.44
1:E:219:MET:O	1:E:220:ILE:HD13	2.17	0.44
1:C:282:LEU:HD22	1:C:284:ILE:HG23	1.98	0.44
1:B:183:ARG:HB2	1:B:201:THR:HA	1.99	0.44
1:F:92:THR:CG2	1:F:104:ALA:HB1	2.47	0.44
1:F:122:ALA:HB2	4:F:395:HOH:O	2.18	0.44
1:F:3:GLN:NE2	1:F:4:SER:O	2.51	0.44
1:B:42:ALA:HB3	1:B:43:PRO:HD3	2.00	0.43
3:E:345:GOL:H12	1:F:139:ASN:HB2	1.97	0.43
1:E:133:PRO:O	1:E:135:VAL:HG23	2.18	0.43
1:F:230:VAL:HG12	1:F:234:SER:HB2	1.99	0.43
1:C:319:ASN:HD22	1:C:322:ASN:H	1.66	0.43
1:C:311:GLN:HB2	1:C:314:LEU:HD11	2.00	0.43
1:F:8:LEU:HB2	1:F:93:LEU:CD2	2.47	0.43
1:D:261:LEU:C	1:D:261:LEU:CD1	2.92	0.43
1:D:268:ILE:HG13	1:D:268:ILE:O	2.19	0.43
1:E:14:THR:HG21	4:E:350:HOH:O	2.18	0.43
1:E:201:THR:HG23	1:E:219:MET:SD	2.58	0.43
1:E:8:LEU:HD12	1:E:9:ALA:N	2.33	0.43
1:E:41:VAL:HG12	1:E:45:LEU:CD1	2.49	0.43
1:E:277:ILE:O	1:E:315:LEU:HD12	2.19	0.43
1:A:102:SER:CA	1:A:131:LEU:HD22	2.49	0.43
1:D:282:LEU:HD22	1:D:284:ILE:HG23	2.01	0.43
1:E:154:GLU:HA	1:E:157:LEU:HD12	2.01	0.43
1:D:153:LEU:HG	1:F:27:ALA:HB2	2.00	0.42
1:E:261:LEU:HD12	1:E:277:ILE:HG23	2.01	0.42
1:A:27:ALA:O	1:A:28:LEU:HD23	2.20	0.42
1:D:311:GLN:HB2	1:D:314:LEU:CD1	2.44	0.42
1:B:268:ILE:HG21	1:B:272:LEU:CD1	2.36	0.42
1:C:44:ILE:HD13	1:C:64:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:47:TYR:OH	1:D:95:ALA:HB2	2.20	0.42
1:F:82:LEU:O	1:F:89:LEU:HD11	2.20	0.42
1:C:26:TYR:HB2	1:C:111:HIS:CD2	2.55	0.42
1:F:56:PHE:CE2	1:F:91:ALA:HB2	2.55	0.42
1:A:27:ALA:HB2	1:B:153:LEU:HD13	2.02	0.42
1:C:41:VAL:HG13	1:C:64:LEU:HD22	2.01	0.42
1:D:8:LEU:HD13	1:D:93:LEU:HD22	2.00	0.42
1:D:185:ALA:HB1	1:D:204:HIS:CE1	2.55	0.42
1:E:187:THR:HG23	1:F:166:SER:HA	2.01	0.42
1:C:16:ASN:OD1	1:C:20:ALA:HB3	2.19	0.42
1:D:183:ARG:HD2	1:E:168:ASP:OD1	2.20	0.42
1:F:179:PRO:O	1:F:182:VAL:HG22	2.20	0.42
1:C:8:LEU:HD23	1:C:33:PRO:CG	2.48	0.42
1:C:16:ASN:CG	1:C:20:ALA:HB3	2.45	0.42
1:F:56:PHE:HB2	1:F:60:GLN:HB2	2.01	0.42
1:A:8:LEU:C	1:A:8:LEU:HD12	2.45	0.41
1:A:42:ALA:HB3	1:A:43:PRO:HD3	2.01	0.41
1:E:66:ASP:CG	4:E:1004:HOH:O	2.63	0.41
1:B:16:ASN:HB3	1:B:22:LEU:HD21	2.02	0.41
1:C:90:VAL:HG22	1:C:92:THR:CG2	2.50	0.41
1:F:53:ALA:CB	1:F:104:ALA:HB3	2.50	0.41
1:A:153:LEU:HD22	1:A:157:LEU:CD1	2.51	0.41
2:A:345:P0A:HA	1:B:209:ASN:ND2	2.34	0.41
1:C:108:LEU:HD12	1:C:108:LEU:C	2.45	0.41
1:E:108:LEU:O	1:E:108:LEU:HD12	2.20	0.41
1:F:128:ILE:HG22	4:F:460:HOH:O	2.20	0.41
1:A:189:ARG:HD3	1:A:206:GLY:O	2.20	0.41
1:F:116:ARG:NH2	1:F:179:PRO:HB2	2.35	0.41
1:A:139:ASN:CB	3:A:346:GOL:H12	2.49	0.41
1:A:268:ILE:HG21	1:A:272:LEU:CD1	2.51	0.41
1:A:282:LEU:HD12	1:A:317:ARG:HA	2.03	0.41
1:E:201:THR:CG2	1:E:219:MET:HA	2.49	0.41
2:A:351:P0A:HAH	1:C:203:MET:SD	2.61	0.41
1:B:81:ARG:HB3	1:C:157:LEU:HD11	2.03	0.41
1:C:12:VAL:O	1:C:12:VAL:HG13	2.20	0.41
1:D:115:HIS:HD2	4:D:939:HOH:O	2.04	0.41
1:F:153:LEU:C	1:F:153:LEU:CD2	2.94	0.41
1:C:9:ALA:HB1	1:C:26:TYR:CD2	2.57	0.40
1:A:145:LEU:HD23	1:A:145:LEU:HA	1.94	0.40
1:D:252:ILE:C	1:D:252:ILE:HD12	2.46	0.40
1:E:69:LYS:NZ	4:E:1010:HOH:O	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:17:ARG:NH2	4:C:784:HOH:O	2.54	0.40
1:F:82:LEU:HB3	1:F:89:LEU:HD13	2.03	0.40
1:F:129:PHE:O	1:F:169:LYS:NZ	2.38	0.40
1:D:156:ARG:HG3	1:F:25:PHE:CD2	2.56	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	315/347 (91%)	302 (96%)	13 (4%)	0	100	100
1	B	312/347 (90%)	302 (97%)	10 (3%)	0	100	100
1	C	317/347 (91%)	306 (96%)	11 (4%)	0	100	100
1	D	315/347 (91%)	307 (98%)	8 (2%)	0	100	100
1	E	303/347 (87%)	299 (99%)	4 (1%)	0	100	100
1	F	317/347 (91%)	303 (96%)	14 (4%)	0	100	100
All	All	1879/2082 (90%)	1819 (97%)	60 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	244/263 (93%)	235 (96%)	9 (4%)	30	22
1	B	241/263 (92%)	228 (95%)	13 (5%)	20	12
1	C	244/263 (93%)	233 (96%)	11 (4%)	24	17
1	D	244/263 (93%)	231 (95%)	13 (5%)	20	12
1	E	236/263 (90%)	224 (95%)	12 (5%)	21	13
1	F	244/263 (93%)	231 (95%)	13 (5%)	20	12
All	All	1453/1578 (92%)	1382 (95%)	71 (5%)	22	14

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	93	LEU
1	A	135	VAL
1	A	153	LEU
1	A	182	VAL
1	A	243	MET
1	A	251	ASN
1	A	306	ARG
1	A	319	ASN
1	B	5	LEU
1	B	12	VAL
1	B	41	VAL
1	B	82	LEU
1	B	177	VAL
1	B	182	VAL
1	B	201	THR
1	B	243	MET
1	B	252	ILE
1	B	282	LEU
1	B	297	ASN
1	B	306	ARG
1	B	315	LEU
1	C	8	LEU
1	C	14	THR
1	C	35	SER
1	C	59	GLN
1	C	93	LEU
1	C	153	LEU
1	C	182	VAL
1	C	235	ASP

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Mol	Chain	Res	Type
1	C	252	ILE
1	C	261	LEU
1	C	329	THR
1	D	4	SER
1	D	8	LEU
1	D	35	SER
1	D	93	LEU
1	D	150	GLU
1	D	153	LEU
1	D	261	LEU
1	D	265	ASN
1	D	282	LEU
1	D	284	ILE
1	D	300	VAL
1	D	306	ARG
1	D	319	ASN
1	E	8	LEU
1	E	14	THR
1	E	69	LYS
1	E	92	THR
1	E	93	LEU
1	E	150	GLU
1	E	177	VAL
1	E	198	GLU
1	E	243	MET
1	E	252	ILE
1	E	282	LEU
1	E	329	THR
1	F	8	LEU
1	F	14	THR
1	F	35	SER
1	F	59	GLN
1	F	92	THR
1	F	96	GLU
1	F	150	GLU
1	F	153	LEU
1	F	282	LEU
1	F	284	ILE
1	F	306	ARG
1	F	319	ASN
1	F	322	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	18	GLN
1	A	52	GLN
1	A	59	GLN
1	A	60	GLN
1	A	115	HIS
1	A	139	ASN
1	A	251	ASN
1	A	276	ASN
1	A	297	ASN
1	A	319	ASN
1	B	3	GLN
1	B	18	GLN
1	B	52	GLN
1	B	60	GLN
1	B	75	GLN
1	B	115	HIS
1	B	134	ASN
1	C	15	GLN
1	C	52	GLN
1	C	59	GLN
1	C	60	GLN
1	C	63	GLN
1	C	121	HIS
1	C	265	ASN
1	C	276	ASN
1	C	297	ASN
1	C	319	ASN
1	C	321	GLN
1	D	15	GLN
1	D	52	GLN
1	D	60	GLN
1	D	86	GLN
1	D	115	HIS
1	D	134	ASN
1	D	139	ASN
1	D	319	ASN
1	E	3	GLN
1	E	15	GLN
1	E	18	GLN
1	E	52	GLN
1	E	59	GLN

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Mol	Chain	Res	Type
1	E	60	GLN
1	E	121	HIS
1	E	134	ASN
1	E	251	ASN
1	F	60	GLN
1	F	134	ASN
1	F	276	ASN
1	F	319	ASN

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](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	P0A	C	351	-	10,11,11	1.01	0	10,13,13	0.91	0
2	P0A	E	351	-	10,11,11	1.08	0	10,13,13	0.90	1 (10%)
2	P0A	D	351	-	10,11,11	1.07	0	10,13,13	0.70	0
2	P0A	A	351	-	10,11,11	1.13	0	10,13,13	0.82	0
3	GOL	A	346	-	5,5,5	0.34	0	5,5,5	1.72	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	P0A	F	351	-	10,11,11	1.06	0	10,13,13	0.90	0
3	GOL	E	345	-	5,5,5	0.39	0	5,5,5	1.07	0
2	P0A	A	345	-	10,11,11	1.09	0	10,13,13	1.01	1 (10%)

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	P0A	C	351	-	-	4/11/11/11	-
2	P0A	E	351	-	-	5/11/11/11	-
2	P0A	D	351	-	-	2/11/11/11	-
2	P0A	A	351	-	-	5/11/11/11	-
3	GOL	A	346	-	-	1/4/4/4	-
2	P0A	F	351	-	-	2/11/11/11	-
3	GOL	E	345	-	-	1/4/4/4	-
2	P0A	A	345	-	-	6/11/11/11	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	346	GOL	O2-C2-C1	2.70	120.35	109.18
2	A	345	P0A	OAB-CAJ-CAH	-2.26	115.91	123.09
2	E	351	P0A	OAB-CAJ-CAH	-2.15	116.27	123.09

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	351	P0A	O-C-CA-N
2	A	345	P0A	O-C-CA-N
2	A	351	P0A	CAG-CAF-CAH-CAJ
2	A	345	P0A	OXT-C-CA-N
2	E	351	P0A	CAH-CAF-CAG-CB
2	A	345	P0A	CAH-CAF-CAG-CB
2	C	351	P0A	CAH-CAF-CAG-CB
2	A	345	P0A	CAF-CAG-CB-CA

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Mol	Chain	Res	Type	Atoms
2	E	351	P0A	CAF-CAG-CB-CA
3	A	346	GOL	O2-C2-C3-O3
2	D	351	P0A	CAH-CAF-CAG-CB
2	A	351	P0A	OXT-C-CA-N
2	F	351	P0A	C-CA-CB-CAG
2	A	351	P0A	CAH-CAF-CAG-CB
2	A	345	P0A	O-C-CA-CB
2	D	351	P0A	CAG-CAF-CAH-CAJ
2	A	345	P0A	OXT-C-CA-CB
2	E	351	P0A	CAG-CAF-CAH-CAJ
2	E	351	P0A	CAF-CAH-CAJ-OAD
2	E	351	P0A	CAF-CAH-CAJ-OAB
2	C	351	P0A	CAF-CAH-CAJ-OAB
2	C	351	P0A	CAG-CAF-CAH-CAJ
2	C	351	P0A	CAF-CAH-CAJ-OAD
3	E	345	GOL	O1-C1-C2-O2
2	A	351	P0A	CAF-CAH-CAJ-OAB
2	F	351	P0A	CAF-CAH-CAJ-OAB

There are no ring outliers.

7 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	351	P0A	2	0
2	E	351	P0A	2	0
2	D	351	P0A	1	0
2	A	351	P0A	1	0
3	A	346	GOL	2	0
3	E	345	GOL	5	0
2	A	345	P0A	3	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 ⓘ

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	321/347 (92%)	0.50	29 (9%) 15 16	17, 28, 51, 57	0
1	B	318/347 (91%)	0.49	35 (11%) 10 11	17, 27, 51, 57	0
1	C	321/347 (92%)	0.55	29 (9%) 15 16	18, 28, 51, 57	0
1	D	321/347 (92%)	0.46	28 (8%) 16 17	17, 27, 51, 56	0
1	E	311/347 (89%)	0.48	27 (8%) 16 17	17, 26, 50, 56	0
1	F	321/347 (92%)	0.57	25 (7%) 19 20	18, 28, 50, 56	0
All	All	1913/2082 (91%)	0.51	173 (9%) 15 16	17, 27, 51, 57	0

All (173) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	325	VAL	5.5
1	E	299	LEU	5.4
1	A	300	VAL	5.3
1	B	280	ALA	5.1
1	D	283	TYR	5.0
1	D	290	VAL	4.9
1	E	324	ALA	4.9
1	E	252	ILE	4.9
1	A	283	TYR	4.8
1	A	290	VAL	4.6
1	F	293	LEU	4.4
1	F	283	TYR	4.3
1	E	305	ALA	4.3
1	E	297	ASN	4.2
1	C	252	ILE	4.2
1	E	280	ALA	4.1
1	F	70	GLY	4.0
1	A	299	LEU	4.0
1	D	299	LEU	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	252	ILE	4.0
1	B	325	VAL	4.0
1	C	283	TYR	4.0
1	D	300	VAL	3.9
1	D	291	ALA	3.8
1	B	300	VAL	3.8
1	F	292	LEU	3.8
1	F	244	GLY	3.7
1	B	283	TYR	3.7
1	B	291	ALA	3.7
1	C	40	ALA	3.7
1	D	323	GLY	3.6
1	B	308	LEU	3.6
1	A	292	LEU	3.6
1	D	302	VAL	3.5
1	E	302	VAL	3.5
1	A	291	ALA	3.5
1	C	292	LEU	3.4
1	B	299	LEU	3.4
1	E	303	VAL	3.3
1	C	293	LEU	3.3
1	E	308	LEU	3.3
1	F	315	LEU	3.3
1	D	298	ALA	3.3
1	F	299	LEU	3.3
1	A	319	ASN	3.3
1	A	245	THR	3.3
1	D	321	GLN	3.2
1	E	283	TYR	3.2
1	C	66	ASP	3.2
1	B	285	THR	3.2
1	A	315	LEU	3.2
1	E	300	VAL	3.1
1	B	298	ALA	3.1
1	D	252	ILE	3.1
1	C	302	VAL	3.1
1	F	252	ILE	3.0
1	C	315	LEU	3.0
1	E	294	ASP	3.0
1	A	318	ARG	3.0
1	B	315	LEU	2.9
1	E	3	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	312	PRO	2.9
1	E	279	GLU	2.9
1	B	305	ALA	2.9
1	A	302	VAL	2.9
1	A	305	ALA	2.9
1	A	321	GLN	2.8
1	F	66	ASP	2.8
1	E	242	THR	2.8
1	D	253	VAL	2.8
1	D	319	ASN	2.8
1	B	302	VAL	2.8
1	C	300	VAL	2.8
1	F	320	SER	2.8
1	D	293	LEU	2.8
1	B	324	ALA	2.7
1	E	298	ALA	2.7
1	A	69	LYS	2.7
1	C	61	ALA	2.7
1	D	254	ILE	2.7
1	C	3	GLN	2.7
1	A	320	SER	2.7
1	B	270	ILE	2.6
1	A	323	GLY	2.6
1	E	290	VAL	2.6
1	F	298	ALA	2.6
1	A	293	LEU	2.6
1	A	308	LEU	2.6
1	B	3	GLN	2.6
1	B	294	ASP	2.6
1	A	325	VAL	2.6
1	D	277	ILE	2.5
1	A	303	VAL	2.5
1	B	303	VAL	2.5
1	B	86	GLN	2.5
1	D	284	ILE	2.5
1	A	84	GLU	2.5
1	C	57	THR	2.5
1	D	325	VAL	2.5
1	F	302	VAL	2.5
1	D	324	ALA	2.5
1	C	253	VAL	2.5
1	F	254	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	245	THR	2.4
1	F	57	THR	2.4
1	B	293	LEU	2.4
1	C	299	LEU	2.4
1	B	251	ASN	2.4
1	B	327	CYS	2.4
1	D	292	LEU	2.4
1	A	324	ALA	2.4
1	B	309	ALA	2.4
1	C	297	ASN	2.4
1	F	278	VAL	2.4
1	B	279	GLU	2.3
1	D	305	ALA	2.3
1	C	19	GLU	2.3
1	C	296	GLN	2.3
1	B	284	ILE	2.3
1	B	312	PRO	2.3
1	D	243	MET	2.3
1	A	304	LYS	2.3
1	E	301	LYS	2.3
1	B	253	VAL	2.3
1	A	282	LEU	2.3
1	F	314	LEU	2.3
1	B	301	LYS	2.3
1	B	268	ILE	2.3
1	F	50	GLY	2.3
1	C	312	PRO	2.2
1	C	282	LEU	2.2
1	D	315	LEU	2.2
1	A	4	SER	2.2
1	A	298	ALA	2.2
1	D	66	ASP	2.2
1	E	309	ALA	2.2
1	A	252	ILE	2.2
1	C	70	GLY	2.2
1	F	323	GLY	2.2
1	B	297	ASN	2.2
1	D	251	ASN	2.2
1	E	304	LYS	2.2
1	A	244	GLY	2.2
1	B	292	LEU	2.2
1	B	286	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	298	ALA	2.2
1	F	305	ALA	2.2
1	E	268	ILE	2.1
1	F	58	SER	2.1
1	B	310	GLY	2.1
1	F	262	ILE	2.1
1	D	244	GLY	2.1
1	D	303	VAL	2.1
1	E	278	VAL	2.1
1	C	314	LEU	2.1
1	E	282	LEU	2.1
1	C	67	ALA	2.1
1	C	284	ILE	2.1
1	A	251	ASN	2.1
1	D	3	GLN	2.1
1	F	59	GLN	2.1
1	F	282	LEU	2.1
1	F	95	ALA	2.1
1	B	304	LYS	2.0
1	C	316	PHE	2.0
1	E	316	PHE	2.0
1	C	242	THR	2.0
1	B	282	LEU	2.0
1	E	315	LEU	2.0
1	F	291	ALA	2.0
1	C	58	SER	2.0
1	C	80	SER	2.0
1	C	322	ASN	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 [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	A	346	6/6	0.77	0.20	18,25,27,27	0
2	P0A	E	351	12/12	0.81	0.11	33,38,38,39	0
2	P0A	A	345	12/12	0.84	0.10	31,36,37,37	0
3	GOL	E	345	6/6	0.88	0.18	25,27,28,28	0
2	P0A	C	351	12/12	0.89	0.10	40,43,44,44	0
2	P0A	F	351	12/12	0.89	0.11	42,43,43,43	0
2	P0A	D	351	12/12	0.93	0.07	20,27,29,30	0
2	P0A	A	351	12/12	0.95	0.07	20,29,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.