



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:48 AM GMT

PDB ID : 2YEP
Title : STRUCTURE OF AN N-TERMINAL NUCLEOPHILE (NTN) HYDROLASE, OAT2, IN COMPLEX WITH GLUTAMATE
Authors : Chowdhury, R.; Iqbal, A.; Clifton, I.J.; Schofield, C.J.
Deposited on : 2011-03-29
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

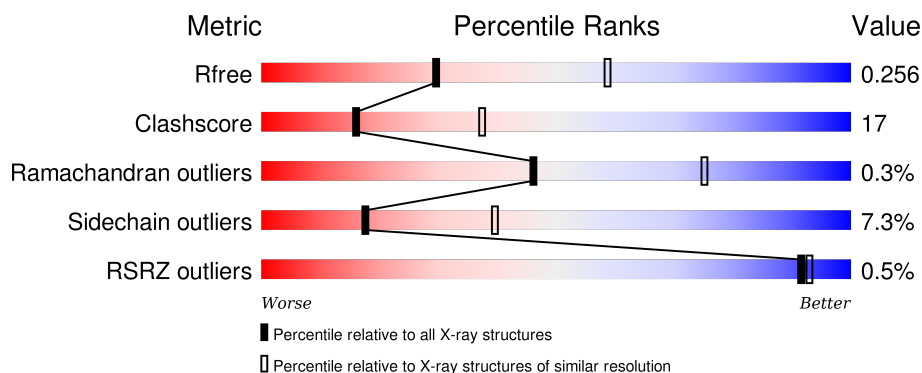
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	180	<div> <div>59%</div> <div>32%</div> <div>• •</div> </div>
1	C	180	<div> <div>2%</div> <div>69%</div> <div>23%</div> <div>• •</div> </div>
1	E	180	<div> <div>58%</div> <div>35%</div> <div>• •</div> </div>
1	G	180	<div> <div>2%</div> <div>63%</div> <div>30%</div> <div>• •</div> </div>
2	B	213	<div> <div>67%</div> <div>29%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	213	
2	H	213	
3	F	213	

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	TH5	B	181	-	-	X	-
4	GLU	B	1601	-	-	X	X
4	GLU	C	1602	-	-	X	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11651 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 GLUTAMATE N-ACETYLTRANSFERASE 2 ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	173	Total	C	N	O	S	0	0	0
			1266	783	234	243	6			
1	C	173	Total	C	N	O	S	0	0	0
			1266	783	234	243	6			
1	E	173	Total	C	N	O	S	0	0	0
			1266	783	234	243	6			
1	G	173	Total	C	N	O	S	0	0	0
			1266	783	234	243	6			

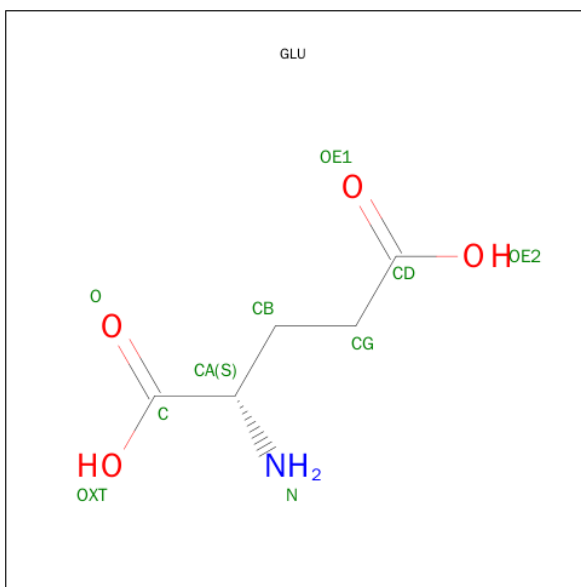
- Molecule 2 is a protein called GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1605	992	279	329	5			
2	D	210	Total	C	N	O	S	0	0	0
			1579	975	276	323	5			
2	H	195	Total	C	N	O	S	0	0	0
			1471	912	253	301	5			

- Molecule 3 is a protein called GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	197	Total	C	N	O	S	0	0	0
			1487	923	255	304	5			

- Molecule 4 is GLUTAMIC ACID (three-letter code: GLU) (formula: C₅H₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	0	0
			9	5	1	3		
4	B	1	Total	C	N	O	0	0
			9	5	1	3		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	C	O	0	0
			4	2	2		
5	G	1	Total	C	O	0	0
			4	2	2		

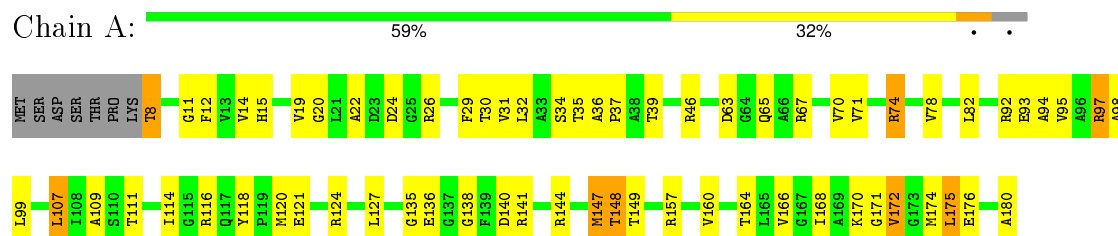
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	71	Total 71	O 71	0	0
6	B	62	Total 62	O 62	0	0
6	C	58	Total 58	O 58	0	0
6	D	55	Total 55	O 55	0	0
6	E	41	Total 41	O 41	0	0
6	F	50	Total 50	O 50	0	0
6	G	43	Total 43	O 43	0	0
6	H	39	Total 39	O 39	0	0

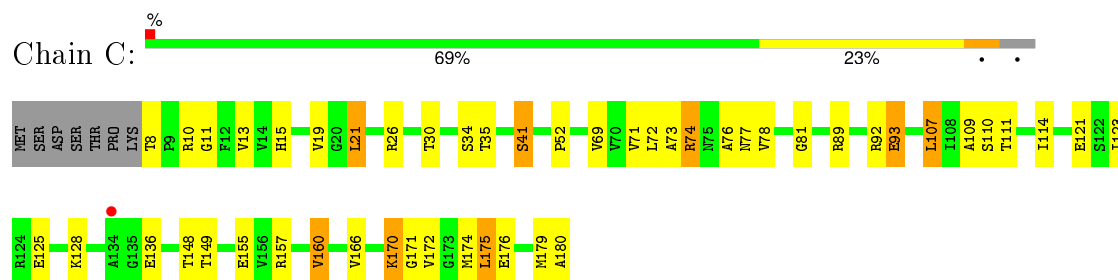
3 Residue-property plots

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.

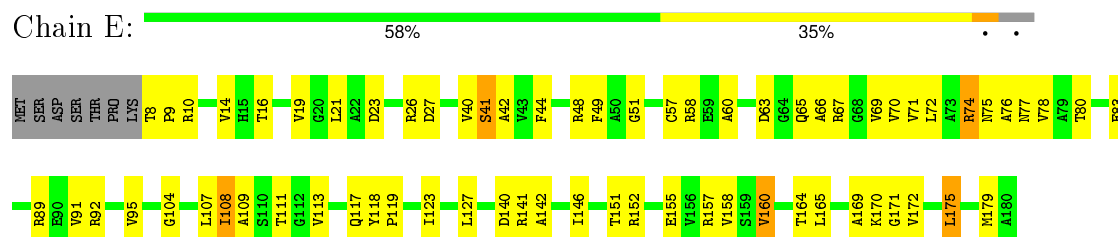
• Molecule 1: GLUTAMATE N-ACETYLTRANSFERASE 2 ALPHA CHAIN



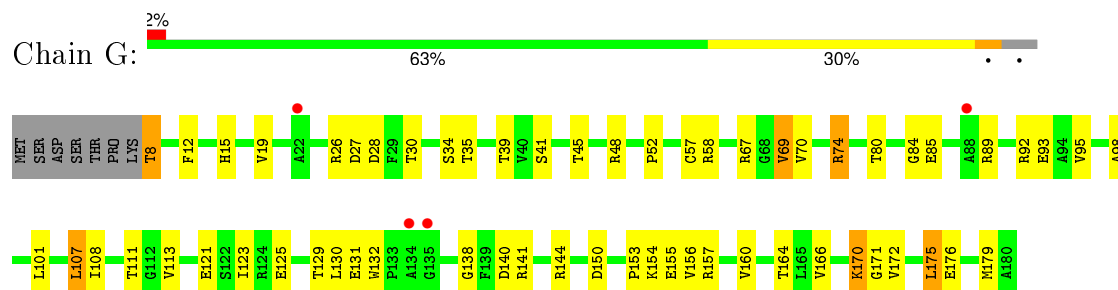
• Molecule 1: GLUTAMATE N-ACETYLTRANSFERASE 2 ALPHA CHAIN



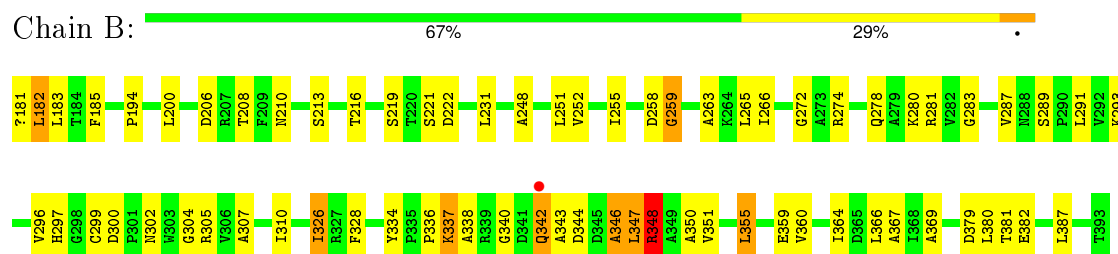
• Molecule 1: GLUTAMATE N-ACETYLTRANSFERASE 2 ALPHA CHAIN



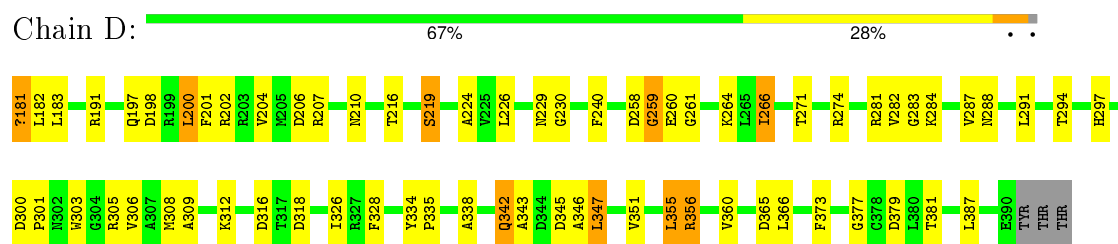
• Molecule 1: GLUTAMATE N-ACETYLTRANSFERASE 2 ALPHA CHAIN



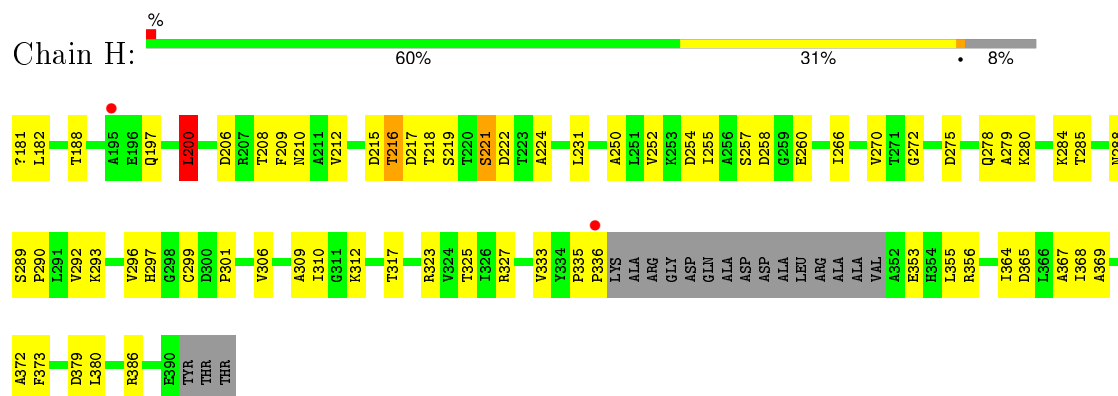
- Molecule 2: GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN



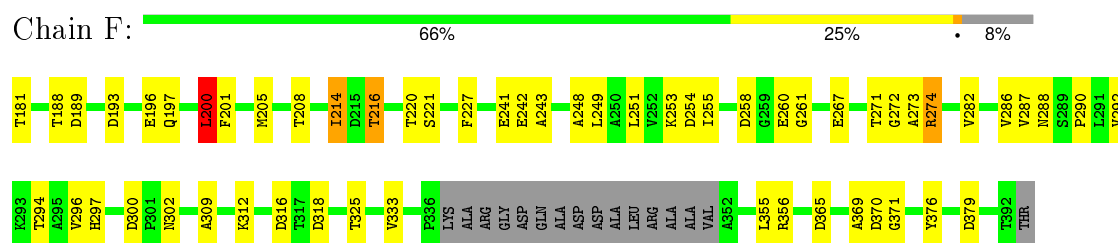
- Molecule 2: GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN



- Molecule 2: GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN



- Molecule 3: GLUTAMATE N-ACETYLTRANSFERASE 2 BETA CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.22Å 73.42Å 172.33Å 90.00° 93.26° 90.00°	Depositor
Resolution (Å)	61.12 – 2.70 67.53 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.0 (61.12-2.70) 96.1 (67.53-2.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.65 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.236 0.214 , 0.256	Depositor DCC
R_{free} test set	2059 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	14.4	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	2 of 40640 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	11651	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.72% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: TH5, ACT

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	0.75	0/1284	0.92	3/1742 (0.2%)
1	C	0.75	0/1284	0.85	2/1742 (0.1%)
1	E	0.68	0/1284	0.82	1/1742 (0.1%)
1	G	0.68	0/1284	0.80	0/1742
2	B	0.77	1/1616 (0.1%)	0.90	6/2195 (0.3%)
2	D	0.75	0/1589	0.85	1/2157 (0.0%)
2	H	0.68	0/1480	0.82	1/2009 (0.0%)
3	F	0.71	0/1507	0.84	3/2047 (0.1%)
All	All	0.72	1/11328 (0.0%)	0.85	17/15376 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	299	CYS	CB-SG	-5.69	1.72	1.81

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	MET	CB-CA-C	-8.68	93.03	110.40
2	H	200	LEU	CA-CB-CG	8.56	134.99	115.30
1	A	148	THR	N-CA-C	6.89	129.59	111.00
3	F	200	LEU	CA-CB-CG	6.59	130.47	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	160	VAL	CB-CA-C	-6.48	99.09	111.40
2	B	366	LEU	CB-CA-C	6.23	122.04	110.20
1	C	107	LEU	CA-CB-CG	6.13	129.41	115.30
2	B	348	ARG	N-CA-C	-5.91	95.04	111.00
3	F	370	ASP	N-CA-C	5.84	126.77	111.00
3	F	274	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	E	160	VAL	CB-CA-C	-5.75	100.47	111.40
2	B	346	ALA	CB-CA-C	-5.70	101.55	110.10
2	B	367	ALA	N-CA-CB	5.69	118.06	110.10
2	B	200	LEU	CA-CB-CG	5.62	128.22	115.30
2	B	274	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	107	LEU	CA-CB-CG	5.24	127.36	115.30
2	D	356	ARG	NE-CZ-NH1	5.20	122.90	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	181	TH5	Mainchain

5.2 Too-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 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	1266	0	1274	62	1
1	C	1266	0	1274	41	0
1	E	1266	0	1274	53	0
1	G	1266	0	1274	55	0
2	B	1605	0	1559	63	0
2	D	1579	0	1536	57	0
2	H	1471	0	1428	54	0
3	F	1487	0	1440	42	0
4	B	9	0	5	7	0
4	C	9	0	5	5	0
5	E	4	0	3	0	0
5	G	4	0	3	0	0
6	A	71	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	62	0	0	7	1
6	C	58	0	0	4	0
6	D	55	0	0	5	0
6	E	41	0	0	8	0
6	F	50	0	0	2	0
6	G	43	0	0	9	0
6	H	39	0	0	3	0
All	All	11651	0	11075	377	1

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 17.

All (377) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:ASP:HB3	6:A:2054:HOH:O	1.45	1.14
1:G:153:PRO:HB2	6:G:2033:HOH:O	1.46	1.13
1:E:60:ALA:O	6:E:2019:HOH:O	1.69	1.09
1:A:26:ARG:NH2	1:A:147:MET:O	1.93	1.01
1:G:74:ARG:NH1	1:G:74:ARG:HB2	1.77	0.99
1:G:74:ARG:HH11	1:G:74:ARG:HB2	1.31	0.92
2:B:347:LEU:C	2:B:347:LEU:HD12	1.89	0.92
1:G:175:LEU:HA	6:G:2040:HOH:O	1.68	0.91
1:A:8:THR:OG1	1:A:157:ARG:NH1	2.03	0.90
2:B:346:ALA:O	2:B:347:LEU:HG	1.73	0.88
1:G:132:TRP:O	6:G:2031:HOH:O	1.93	0.85
6:G:2043:HOH:O	2:H:222:ASP:OD2	1.93	0.85
3:F:282:VAL:O	3:F:286:VAL:HG23	1.76	0.85
2:B:263:ALA:CB	6:B:2028:HOH:O	2.24	0.84
1:G:140:ASP:O	1:G:144:ARG:HG3	1.78	0.83
1:C:74:ARG:HH11	1:C:74:ARG:HB2	1.43	0.83
1:C:19:VAL:HG11	1:C:123:ILE:HD13	1.64	0.80
2:H:310:ILE:HD11	2:H:364:ILE:HD13	1.65	0.79
2:H:306:VAL:O	2:H:310:ILE:HG12	1.84	0.78
1:C:19:VAL:O	1:C:19:VAL:HG12	1.85	0.77
1:A:171:GLY:HA2	4:B:1601:GLU:OE2	1.85	0.76
2:H:336:PRO:HA	6:H:2038:HOH:O	1.84	0.76
1:E:158:VAL:HG11	3:F:243:ALA:HB1	1.66	0.76
2:H:272:GLY:HA2	2:H:369:ALA:O	1.85	0.75
1:A:20:GLY:O	6:A:2004:HOH:O	2.03	0.75
3:F:197:GLN:HA	3:F:200:LEU:HD22	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:VAL:N	4:B:1601:GLU:OE2	2.21	0.74
1:G:172:VAL:HG23	2:H:260:GLU:HG3	1.71	0.73
1:G:157:ARG:HG2	1:G:166:VAL:HG22	1.70	0.73
1:C:171:GLY:O	1:C:175:LEU:HD13	1.88	0.73
2:H:310:ILE:CD1	2:H:364:ILE:HD13	2.18	0.73
3:F:325:THR:HG21	6:F:2039:HOH:O	1.89	0.73
1:G:45:THR:HB	2:H:222:ASP:HB2	1.70	0.72
2:B:266:ILE:HD11	2:B:296:VAL:HG11	1.70	0.72
3:F:290:PRO:O	3:F:294:THR:HG23	1.89	0.71
2:B:347:LEU:C	2:B:347:LEU:CD1	2.58	0.71
1:A:36:ALA:HB2	1:G:35:THR:O	1.92	0.70
1:C:172:VAL:HG23	2:D:260:GLU:HG3	1.73	0.70
2:B:206:ASP:HA	2:B:210:ASN:HB2	1.73	0.70
1:C:157:ARG:HG2	1:C:166:VAL:HG22	1.74	0.70
1:C:148:THR:OG1	4:C:1602:GLU:OE1	2.07	0.69
1:G:19:VAL:HG11	1:G:123:ILE:HG13	1.73	0.69
2:B:263:ALA:HB3	6:B:2028:HOH:O	1.90	0.69
2:B:346:ALA:O	2:B:347:LEU:CG	2.40	0.69
1:A:176:GLU:HG3	2:B:213:SER:HB2	1.73	0.69
2:B:208:THR:HB	2:B:252:VAL:HG21	1.74	0.68
2:B:346:ALA:O	2:B:347:LEU:CB	2.41	0.68
2:B:297:HIS:HB2	2:B:379:ASP:HB2	1.75	0.68
1:E:71:VAL:HG22	1:E:109:ALA:HB3	1.74	0.68
2:B:291:LEU:HB3	2:B:305:ARG:HD3	1.75	0.68
2:D:345:ASP:OD2	6:D:2047:HOH:O	2.11	0.68
1:A:144:ARG:HD2	6:A:2056:HOH:O	1.94	0.68
2:H:280:LYS:HD2	2:H:373:PHE:CZ	2.29	0.68
1:E:119:PRO:O	1:E:123:ILE:HG13	1.95	0.67
2:B:343:ALA:O	2:B:347:LEU:HG	1.93	0.67
2:H:335:PRO:HB2	2:H:336:PRO:HD2	1.76	0.65
2:D:343:ALA:O	2:D:347:LEU:HD22	1.96	0.65
1:E:71:VAL:HA	1:E:109:ALA:O	1.95	0.65
1:E:91:VAL:O	1:E:95:VAL:HG23	1.97	0.65
2:B:347:LEU:HD12	2:B:348:ARG:N	2.12	0.64
1:A:140:ASP:CB	6:A:2054:HOH:O	2.22	0.64
3:F:356:ARG:NH1	2:H:323:ARG:HD2	2.12	0.63
2:B:326:ILE:CG1	2:B:334:TYR:HB3	2.29	0.63
2:H:181:TH5:HG22	2:H:221:SER:HB2	1.81	0.62
2:D:283:GLY:HA3	6:D:2021:HOH:O	1.98	0.62
2:D:207:ARG:HH12	2:D:373:PHE:HA	1.65	0.62
1:C:111:THR:HB	2:D:183:LEU:HD11	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:26:ARG:HD2	1:G:27:ASP:O	2.00	0.62
2:B:248:ALA:O	2:B:252:VAL:HG23	2.00	0.62
2:D:291:LEU:HB3	2:D:305:ARG:HD3	1.82	0.62
2:D:274:ARG:HD3	2:D:318:ASP:OD2	2.00	0.61
1:C:73:ALA:O	1:C:74:ARG:HB2	1.99	0.61
1:E:10:ARG:HG3	3:F:189:ASP:OD2	2.00	0.61
2:H:280:LYS:HD2	2:H:373:PHE:CE1	2.36	0.60
1:G:121:GLU:HA	1:G:121:GLU:OE1	2.01	0.60
2:B:347:LEU:HA	2:B:350:ALA:HB3	1.83	0.60
2:B:326:ILE:HG12	2:B:334:TYR:HB3	1.83	0.60
1:E:27:ASP:OD2	6:E:2005:HOH:O	2.16	0.60
3:F:356:ARG:HH12	2:H:323:ARG:HD2	1.67	0.60
1:G:67:ARG:NE	1:G:101:LEU:HD22	2.17	0.60
2:D:266:ILE:N	2:D:266:ILE:CD1	2.64	0.60
2:B:231:LEU:HD22	2:H:231:LEU:HD21	1.85	0.59
2:B:342:GLN:OE1	2:B:342:GLN:N	2.35	0.59
3:F:200:LEU:HD23	3:F:201:PHE:N	2.18	0.59
1:G:70:VAL:HG13	1:G:108:ILE:HG13	1.83	0.59
2:D:197:GLN:HA	2:D:200:LEU:HD22	1.84	0.59
1:A:8:THR:HG1	1:A:157:ARG:HH11	1.49	0.58
2:B:263:ALA:O	2:D:274:ARG:HD2	2.02	0.58
2:B:297:HIS:ND1	2:B:381:THR:HG22	2.17	0.58
2:H:208:THR:HB	2:H:252:VAL:HG21	1.85	0.58
6:E:2022:HOH:O	3:F:221:SER:HB3	2.03	0.58
2:D:200:LEU:HD23	2:D:201:PHE:N	2.18	0.58
1:E:23:ASP:O	1:G:48:ARG:NH2	2.37	0.58
4:C:1602:GLU:HG2	2:D:181:TH5:N	2.18	0.58
1:E:40:VAL:HG21	1:E:66:ALA:HB3	1.86	0.57
1:A:97:ARG:HG2	1:A:98:ALA:N	2.19	0.57
2:B:258:ASP:O	2:B:259:GLY:C	2.42	0.57
2:H:266:ILE:HG12	2:H:296:VAL:HG11	1.86	0.57
2:H:289:SER:HB3	2:H:292:VAL:HB	1.86	0.57
1:G:131:GLU:HB2	6:G:2030:HOH:O	2.03	0.57
1:C:148:THR:OG1	1:C:149:THR:N	2.38	0.57
1:E:70:VAL:HG13	1:E:108:ILE:HD12	1.85	0.57
3:F:274:ARG:HD3	3:F:318:ASP:OD2	2.03	0.57
2:B:181:TH5:HAA1	2:B:219:SER:HB2	1.87	0.57
1:E:51:GLY:HA3	6:E:2013:HOH:O	2.04	0.57
2:D:297:HIS:HB2	2:D:379:ASP:HB2	1.87	0.57
1:G:140:ASP:HB3	6:G:2035:HOH:O	2.04	0.56
1:G:176:GLU:OE1	2:H:216:THR:HA	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:MET:O	1:A:124:ARG:HG3	2.05	0.56
1:A:74:ARG:HB2	1:A:74:ARG:HH11	1.71	0.56
2:B:310:ILE:HD11	2:B:364:ILE:HD13	1.86	0.56
1:C:19:VAL:CG1	1:C:123:ILE:HD13	2.36	0.56
1:C:19:VAL:O	1:C:19:VAL:CG1	2.54	0.56
1:A:46:ARG:NH2	6:A:2013:HOH:O	2.39	0.56
1:C:89:ARG:O	1:C:93:GLU:HB3	2.05	0.55
2:D:281:ARG:NH1	2:D:316:ASP:OD1	2.39	0.55
2:B:344:ASP:O	2:B:348:ARG:HB2	2.06	0.55
1:A:19:VAL:CG2	1:A:29:PHE:HB2	2.37	0.55
2:B:355:LEU:HD12	2:B:360:VAL:CG2	2.36	0.55
2:H:206:ASP:HA	2:H:210:ASN:HB2	1.87	0.55
3:F:297:HIS:HB2	3:F:379:ASP:HB2	1.89	0.55
1:A:121:GLU:HA	1:A:121:GLU:OE1	2.07	0.55
2:B:293:LYS:HE3	2:B:380:LEU:HB2	1.88	0.55
1:A:78:VAL:HB	1:A:114:ILE:HD12	1.88	0.54
3:F:216:THR:HG22	3:F:288:ASN:OD1	2.07	0.54
2:H:325:THR:OG1	2:H:365:ASP:HB3	2.07	0.54
2:D:387:LEU:O	2:D:387:LEU:HD23	2.08	0.54
1:C:69:VAL:HG23	1:C:107:LEU:HB2	1.89	0.53
1:E:165:LEU:HD12	3:F:188:THR:HB	1.90	0.53
2:H:212:VAL:HB	2:H:252:VAL:HG13	1.89	0.53
1:E:76:ALA:O	1:E:78:VAL:HG23	2.08	0.53
1:A:174:MET:SD	2:B:181:TH5:HAA1	2.47	0.53
2:B:272:GLY:HA2	2:B:369:ALA:O	2.08	0.53
1:E:111:THR:OG1	3:F:181:THR:HB	2.08	0.53
1:E:151:THR:OG1	3:F:258:ASP:OD1	2.19	0.53
2:B:340:GLY:HA3	6:B:2051:HOH:O	2.08	0.53
1:G:8:THR:HB	1:G:164:THR:HG21	1.91	0.53
1:E:140:ASP:HB3	6:E:2034:HOH:O	2.08	0.52
2:D:259:GLY:O	2:D:260:GLU:C	2.48	0.52
1:A:19:VAL:HG21	1:A:29:PHE:HB2	1.91	0.52
2:D:266:ILE:N	2:D:266:ILE:HD13	2.25	0.52
2:B:182:LEU:HD22	2:B:183:LEU:N	2.25	0.52
1:E:26:ARG:O	6:E:2002:HOH:O	2.19	0.52
3:F:249:LEU:HD11	3:F:253:LYS:HE3	1.92	0.51
1:C:21:LEU:HD11	1:C:72:LEU:HD23	1.91	0.51
2:D:342:GLN:CD	2:D:342:GLN:H	2.12	0.51
1:A:168:ILE:HD11	2:B:185:PHE:HD2	1.75	0.51
1:G:12:PHE:HE2	2:H:188:THR:O	1.93	0.51
1:E:175:LEU:HD22	3:F:255:ILE:HG21	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:VAL:HA	1:A:109:ALA:O	2.11	0.51
2:H:197:GLN:HA	2:H:200:LEU:HD13	1.93	0.51
1:E:41:SER:OG	3:F:197:GLN:HG2	2.11	0.51
2:H:310:ILE:HD11	2:H:364:ILE:CD1	2.38	0.51
1:A:174:MET:SD	2:B:181:TH5:CAA	2.99	0.51
1:E:142:ALA:O	1:E:146:ILE:HG23	2.10	0.51
1:A:14:VAL:HG22	1:A:32:LEU:HD12	1.92	0.51
2:H:270:VAL:O	2:H:372:ALA:HA	2.11	0.50
1:A:171:GLY:O	2:B:255:ILE:HD12	2.10	0.50
1:G:130:LEU:HG	1:G:131:GLU:N	2.27	0.50
2:H:216:THR:HG22	2:H:288:ASN:CG	2.31	0.50
2:B:280:LYS:HA	6:B:2024:HOH:O	2.11	0.50
1:A:148:THR:OG1	4:B:1601:GLU:OE1	2.30	0.49
2:H:250:ALA:O	2:H:254:ASP:HB2	2.12	0.49
3:F:208:THR:HG21	3:F:248:ALA:HB3	1.93	0.49
3:F:273:ALA:HB3	3:F:371:GLY:HA3	1.94	0.49
1:C:35:THR:HB	1:E:67:ARG:NH1	2.26	0.49
1:G:171:GLY:C	1:G:175:LEU:HD13	2.32	0.49
1:A:78:VAL:HG12	1:A:114:ILE:HG23	1.93	0.49
1:G:154:LYS:HE2	2:H:255:ILE:HA	1.95	0.49
2:H:215:ASP:N	2:H:215:ASP:OD1	2.43	0.49
1:A:24:ASP:OD2	1:A:74:ARG:NH2	2.46	0.49
1:E:21:LEU:HD22	1:E:75:ASN:HB3	1.95	0.49
1:G:74:ARG:HB2	1:G:74:ARG:CZ	2.42	0.49
1:A:37:PRO:HA	1:A:67:ARG:HD2	1.93	0.49
1:G:70:VAL:HG21	1:G:95:VAL:HG21	1.94	0.49
2:B:182:LEU:O	2:B:222:ASP:HA	2.12	0.49
1:G:138:GLY:HA3	1:G:141:ARG:HH21	1.78	0.49
3:F:271:THR:O	3:F:365:ASP:HA	2.13	0.49
1:E:104:GLY:O	6:E:2019:HOH:O	2.20	0.48
1:E:169:ALA:HB2	3:F:251:LEU:HD13	1.94	0.48
1:G:28:ASP:CB	1:G:74:ARG:HH12	2.26	0.48
2:D:197:GLN:O	2:D:200:LEU:CD2	2.62	0.48
2:B:355:LEU:HD12	2:B:360:VAL:HG21	1.95	0.48
1:G:12:PHE:CE2	2:H:188:THR:O	2.66	0.48
2:H:309:ALA:HA	2:H:312:LYS:HG3	1.95	0.48
1:G:74:ARG:HG3	6:G:2016:HOH:O	2.13	0.48
2:B:181:TH5:N	4:B:1601:GLU:HB2	2.28	0.48
2:H:272:GLY:CA	2:H:369:ALA:O	2.59	0.48
1:E:21:LEU:HD11	1:E:72:LEU:HB3	1.95	0.48
2:D:334:TYR:CD1	2:D:335:PRO:HA	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:VAL:HA	1:C:109:ALA:O	2.12	0.48
2:D:300:ASP:OD1	2:D:301:PRO:HD2	2.13	0.48
1:A:22:ALA:H	1:A:74:ARG:HH12	1.62	0.48
3:F:267:GLU:HG3	3:F:376:TYR:CE2	2.49	0.48
1:C:26:ARG:HB3	1:C:74:ARG:HH21	1.79	0.48
2:B:337:LYS:H	2:B:337:LYS:HD2	1.78	0.48
2:H:327:ARG:HG2	6:H:2033:HOH:O	2.13	0.47
1:E:74:ARG:N	1:E:111:THR:O	2.47	0.47
1:A:149:THR:N	4:B:1601:GLU:OE1	2.35	0.47
2:D:200:LEU:O	2:D:204:VAL:HG23	2.14	0.47
1:E:123:ILE:HG22	1:E:127:LEU:HD11	1.96	0.47
2:H:335:PRO:HG3	6:H:2032:HOH:O	2.14	0.47
2:B:265:LEU:HB3	2:B:359:GLU:HG2	1.95	0.47
2:H:297:HIS:HB2	2:H:379:ASP:HB2	1.96	0.47
1:G:69:VAL:HB	1:G:107:LEU:HB2	1.97	0.47
1:A:12:PHE:O	1:A:136:GLU:HB3	2.15	0.47
1:G:170:LYS:NZ	2:H:181:TH5:O	2.39	0.47
1:C:175:LEU:HA	6:C:2055:HOH:O	2.14	0.47
2:D:346:ALA:HB1	6:D:2048:HOH:O	2.15	0.47
1:A:111:THR:OG1	2:B:181:TH5:HB	2.15	0.47
1:G:12:PHE:CE1	1:G:34:SER:HB2	2.50	0.47
2:B:326:ILE:HG13	2:B:334:TYR:HB3	1.96	0.46
2:D:338:ALA:N	6:D:2045:HOH:O	2.47	0.46
2:H:299:CYS:SG	2:H:356:ARG:HD3	2.54	0.46
1:G:111:THR:HG21	1:G:170:LYS:HE3	1.97	0.46
1:G:111:THR:OG1	1:G:170:LYS:HE3	2.14	0.46
2:B:304:GLY:O	2:B:307:ALA:HB3	2.15	0.46
3:F:272:GLY:HA2	3:F:369:ALA:O	2.15	0.46
1:E:83:GLU:HB3	1:E:117:GLN:HE22	1.80	0.46
2:D:284:LYS:O	2:D:288:ASN:HB2	2.15	0.46
2:H:288:ASN:O	2:H:290:PRO:HD3	2.16	0.46
1:G:52:PRO:HB3	1:G:85:GLU:HG2	1.96	0.46
1:A:171:GLY:C	1:A:175:LEU:CD1	2.84	0.46
1:A:120:MET:O	1:A:124:ARG:CG	2.63	0.46
1:C:15:HIS:O	1:C:30:THR:HA	2.16	0.46
1:A:63:ASP:CG	1:A:65:GLN:HB3	2.35	0.46
2:H:272:GLY:O	2:H:368:ILE:HB	2.16	0.46
1:C:174:MET:HG2	2:D:219:SER:HB2	1.98	0.46
3:F:188:THR:HG22	3:F:227:PHE:O	2.16	0.46
1:C:179:MET:HE1	2:D:224:ALA:N	2.31	0.46
2:H:293:LYS:HE3	2:H:380:LEU:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:8:THR:HB	1:E:164:THR:HG21	1.98	0.46
1:E:60:ALA:HB3	1:E:107:LEU:CD2	2.46	0.45
1:A:171:GLY:C	1:A:175:LEU:HD13	2.36	0.45
1:A:11:GLY:O	1:A:34:SER:HA	2.17	0.45
1:A:170:LYS:HD3	1:A:170:LYS:C	2.37	0.45
1:G:179:MET:HE1	2:H:224:ALA:N	2.30	0.45
1:A:171:GLY:CA	4:B:1601:GLU:OE2	2.59	0.45
4:C:1602:GLU:CG	2:D:181:TH5:N	2.79	0.45
1:C:155:GLU:OE2	1:C:157:ARG:NH1	2.49	0.45
3:F:214:ILE:CD1	3:F:214:ILE:N	2.79	0.45
6:G:2043:HOH:O	2:H:222:ASP:CB	2.64	0.45
1:A:30:THR:O	1:A:70:VAL:HA	2.16	0.45
1:C:13:VAL:HG12	1:C:136:GLU:HA	1.98	0.45
2:H:323:ARG:HB3	2:H:367:ALA:HB3	1.98	0.45
1:E:48:ARG:HA	1:E:48:ARG:NE	2.32	0.45
1:A:15:HIS:HB3	1:A:31:VAL:HB	1.97	0.45
2:B:251:LEU:O	2:B:255:ILE:HG12	2.17	0.45
2:D:305:ARG:O	2:D:308:MET:HB3	2.17	0.45
1:A:39:THR:HG21	2:B:194:PRO:HG3	1.98	0.45
3:F:200:LEU:HD12	3:F:241:GLU:HB2	1.98	0.45
2:D:297:HIS:ND1	2:D:381:THR:HG22	2.32	0.45
2:D:198:ASP:O	2:D:202:ARG:HG3	2.17	0.45
1:C:81:GLY:HA2	6:C:2032:HOH:O	2.17	0.44
2:D:342:GLN:OE1	2:D:342:GLN:N	2.45	0.44
1:G:80:THR:OG1	1:G:84:GLY:HA3	2.17	0.44
1:A:22:ALA:H	1:A:74:ARG:NH1	2.14	0.44
1:A:19:VAL:CG2	1:A:29:PHE:CB	2.96	0.44
1:A:157:ARG:HG2	1:A:166:VAL:HG22	1.99	0.44
1:G:19:VAL:CG1	1:G:123:ILE:HG13	2.43	0.44
1:A:46:ARG:HD3	1:A:180:ALA:HB3	2.00	0.44
1:E:146:ILE:O	1:E:170:LYS:HG3	2.17	0.44
3:F:316:ASP:C	3:F:318:ASP:H	2.21	0.44
2:B:382:GLU:OE1	6:B:2060:HOH:O	2.21	0.44
2:B:297:HIS:CE1	2:B:381:THR:HG22	2.53	0.44
1:E:165:LEU:CD1	3:F:188:THR:HB	2.47	0.44
1:A:15:HIS:O	1:A:30:THR:HA	2.18	0.44
1:A:31:VAL:HG11	1:A:99:LEU:HD11	1.99	0.44
1:A:8:THR:HB	1:A:164:THR:HG21	1.99	0.43
1:A:170:LYS:HZ3	4:B:1601:GLU:HB3	1.81	0.43
2:B:208:THR:HB	2:B:252:VAL:CG2	2.47	0.43
1:E:63:ASP:OD1	1:E:65:GLN:HB3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:THR:HB	1:G:67:ARG:HH12	1.82	0.43
1:G:41:SER:OG	2:H:197:GLN:HG2	2.17	0.43
1:C:15:HIS:HE1	6:C:2044:HOH:O	2.02	0.43
3:F:287:VAL:HG22	6:F:2008:HOH:O	2.18	0.43
1:G:15:HIS:O	1:G:30:THR:HA	2.17	0.43
2:H:284:LYS:O	2:H:285:THR:C	2.53	0.43
1:E:44:PHE:O	1:E:179:MET:HA	2.18	0.43
2:B:300:ASP:O	2:B:302:ASN:N	2.44	0.43
1:A:138:GLY:O	1:A:141:ARG:N	2.51	0.43
1:A:168:ILE:CD1	2:B:185:PHE:HD2	2.30	0.43
1:C:52:PRO:HD2	1:C:77:ASN:O	2.18	0.43
2:B:336:PRO:HB3	6:B:2048:HOH:O	2.18	0.43
1:A:109:ALA:HB1	2:B:183:LEU:HD22	2.00	0.43
1:E:49:PHE:O	3:F:220:THR:HB	2.18	0.43
2:D:306:VAL:HG12	2:D:326:ILE:HD12	2.00	0.43
1:C:69:VAL:HG23	1:C:107:LEU:CB	2.48	0.43
2:D:261:GLY:O	2:D:381:THR:HG21	2.19	0.43
1:A:19:VAL:HG21	1:A:29:PHE:CB	2.48	0.43
1:C:176:GLU:OE1	2:D:216:THR:HA	2.19	0.43
1:A:94:ALA:HB3	1:A:127:LEU:HD13	2.01	0.43
2:B:283:GLY:HA3	6:B:2024:HOH:O	2.18	0.42
1:E:42:ALA:HB1	1:E:44:PHE:CZ	2.54	0.42
1:E:69:VAL:HA	1:E:107:LEU:O	2.18	0.42
1:E:40:VAL:CG1	1:E:41:SER:N	2.82	0.42
3:F:201:PHE:O	3:F:205:MET:HG2	2.20	0.42
2:B:355:LEU:HD12	2:B:360:VAL:HG22	2.01	0.42
3:F:272:GLY:C	3:F:369:ALA:O	2.58	0.42
2:H:217:ASP:O	2:H:218:THR:C	2.58	0.42
1:G:28:ASP:HB2	1:G:74:ARG:HH12	1.83	0.42
1:G:19:VAL:CG1	1:G:19:VAL:O	2.67	0.42
1:G:89:ARG:O	1:G:93:GLU:HB3	2.18	0.42
1:G:140:ASP:O	1:G:144:ARG:CG	2.61	0.42
1:G:170:LYS:HG2	1:G:171:GLY:N	2.34	0.42
1:G:172:VAL:O	2:H:260:GLU:HG3	2.19	0.42
2:D:309:ALA:HA	2:D:312:LYS:HG3	1.99	0.42
1:E:172:VAL:HG23	3:F:260:GLU:HG3	2.01	0.42
2:B:181:TH5:HG22	2:B:221:SER:HB2	2.00	0.42
2:H:208:THR:OG1	2:H:209:PHE:N	2.53	0.42
2:D:355:LEU:HD12	2:D:360:VAL:HG11	2.01	0.42
2:D:191:ARG:HB2	2:D:230:GLY:HA3	2.02	0.42
2:H:278:GLN:O	2:H:279:ALA:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:GLY:O	1:C:34:SER:HA	2.19	0.42
2:B:342:GLN:CD	2:B:342:GLN:N	2.73	0.42
1:E:171:GLY:O	3:F:255:ILE:HG23	2.20	0.42
1:C:76:ALA:O	1:C:110:SER:HB3	2.20	0.42
1:E:8:THR:N	1:E:9:PRO:CD	2.83	0.42
1:G:155:GLU:HG2	1:G:156:VAL:N	2.35	0.42
1:A:26:ARG:HB3	1:A:74:ARG:NH2	2.35	0.42
1:A:172:VAL:HG11	2:B:258:ASP:HB3	2.02	0.42
3:F:292:VAL:O	3:F:296:VAL:HG23	2.19	0.42
2:D:282:VAL:HG21	2:D:366:LEU:HD11	2.02	0.42
1:C:170:LYS:HZ3	4:C:1602:GLU:CD	2.23	0.42
1:E:155:GLU:OE1	1:E:157:ARG:NH1	2.53	0.42
2:D:229:ASN:O	2:D:229:ASN:OD1	2.38	0.42
1:C:125:GLU:O	1:C:128:LYS:HB2	2.20	0.42
2:D:200:LEU:HD11	2:D:240:PHE:HD2	1.85	0.42
2:H:289:SER:HB3	2:H:292:VAL:CG2	2.50	0.42
1:A:70:VAL:HG21	1:A:95:VAL:HG21	2.01	0.41
2:B:278:GLN:HA	2:B:281:ARG:NH1	2.35	0.41
1:C:26:ARG:HB3	1:C:74:ARG:NH2	2.35	0.41
1:C:170:LYS:NZ	2:D:181:TH5:O	2.40	0.41
2:D:379:ASP:N	2:D:379:ASP:OD1	2.53	0.41
2:B:328:PHE:HB2	2:B:351:VAL:HG22	2.01	0.41
1:G:67:ARG:CZ	1:G:101:LEU:HD22	2.51	0.41
1:E:14:VAL:HG11	1:E:142:ALA:HB2	2.01	0.41
1:E:77:ASN:HB2	1:E:118:TYR:CE1	2.56	0.41
2:D:347:LEU:HA	2:D:347:LEU:HD12	1.93	0.41
2:D:197:GLN:O	2:D:200:LEU:HD23	2.19	0.41
2:B:337:LYS:O	2:B:338:ALA:HB3	2.20	0.41
2:D:271:THR:O	2:D:365:ASP:HA	2.21	0.41
1:A:176:GLU:OE1	2:B:216:THR:HA	2.19	0.41
1:E:19:VAL:CG1	1:E:123:ILE:HG12	2.50	0.41
1:E:21:LEU:CD1	1:E:72:LEU:HB3	2.50	0.41
1:C:8:THR:O	1:C:8:THR:HG22	2.20	0.41
2:D:206:ASP:HA	2:D:210:ASN:HB2	2.03	0.41
2:H:301:PRO:HB3	2:H:355:LEU:HD13	2.03	0.41
1:E:158:VAL:CG1	3:F:243:ALA:HB1	2.43	0.41
2:D:258:ASP:O	2:D:259:GLY:C	2.59	0.41
1:G:138:GLY:O	1:G:141:ARG:N	2.49	0.41
1:G:57:CYS:O	1:G:58:ARG:C	2.59	0.41
3:F:300:ASP:O	3:F:302:ASN:N	2.52	0.41
1:A:135:GLY:HA2	6:A:2052:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:170:LYS:HG2	1:C:171:GLY:N	2.36	0.41
2:H:258:ASP:O	2:H:258:ASP:CG	2.58	0.41
1:G:144:ARG:NH1	1:G:153:PRO:HG3	2.36	0.40
1:C:41:SER:OG	2:D:197:GLN:HG2	2.21	0.40
2:D:264:LYS:HG3	6:D:2016:HOH:O	2.21	0.40
3:F:309:ALA:O	3:F:312:LYS:HG3	2.21	0.40
2:D:287:VAL:CG2	2:D:377:GLY:HA3	2.51	0.40
1:E:57:CYS:HA	1:E:107:LEU:CD2	2.51	0.40
1:E:60:ALA:HB3	1:E:107:LEU:HD23	2.02	0.40
2:H:335:PRO:HB2	2:H:336:PRO:CD	2.44	0.40
1:C:157:ARG:NH2	6:C:2047:HOH:O	2.55	0.40
1:G:98:ALA:HB1	1:G:130:LEU:O	2.21	0.40
1:C:180:ALA:O	2:D:202:ARG:NH2	2.52	0.40
1:E:16:THR:HG21	1:E:141:ARG:HB3	2.04	0.40
3:F:309:ALA:HA	3:F:312:LYS:HG3	2.02	0.40
2:D:328:PHE:HB2	2:D:351:VAL:HG22	2.03	0.40
1:E:58:ARG:NE	6:E:2016:HOH:O	2.43	0.40
1:G:74:ARG:HA	6:G:2016:HOH:O	2.20	0.40
4:C:1602:GLU:N	2:D:181:TH5:HN1	2.20	0.40
2:D:343:ALA:O	2:D:347:LEU:CD2	2.67	0.40
2:D:303:TRP:O	2:D:306:VAL:HB	2.21	0.40
3:F:193:ASP:HB3	3:F:196:GLU:HB2	2.03	0.40
1:C:78:VAL:HB	1:C:114:ILE:HD12	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:ARG:NH1	6:B:2060:HOH:O[2_544]	1.88	0.32

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	171/180 (95%)	159 (93%)	11 (6%)	1 (1%)	30	59
1	C	171/180 (95%)	160 (94%)	11 (6%)	0	100	100
1	E	171/180 (95%)	163 (95%)	8 (5%)	0	100	100
1	G	171/180 (95%)	155 (91%)	16 (9%)	0	100	100
2	B	211/213 (99%)	196 (93%)	14 (7%)	1 (0%)	34	63
2	D	208/213 (98%)	192 (92%)	15 (7%)	1 (0%)	34	63
2	H	191/213 (90%)	176 (92%)	15 (8%)	0	100	100
3	F	193/213 (91%)	181 (94%)	11 (6%)	1 (0%)	34	63
All	All	1487/1572 (95%)	1382 (93%)	101 (7%)	4 (0%)	46	75

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	259	GLY
1	A	172	VAL
2	D	259	GLY
3	F	261	GLY

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	128/135 (95%)	117 (91%)	11 (9%)	13	29
1	C	128/135 (95%)	118 (92%)	10 (8%)	16	35
1	E	128/135 (95%)	118 (92%)	10 (8%)	16	35
1	G	128/135 (95%)	115 (90%)	13 (10%)	9	21
2	B	165/165 (100%)	155 (94%)	10 (6%)	23	49
2	D	162/165 (98%)	152 (94%)	10 (6%)	23	49
2	H	153/165 (93%)	142 (93%)	11 (7%)	18	41
3	F	156/166 (94%)	149 (96%)	7 (4%)	34	65
All	All	1148/1201 (96%)	1066 (93%)	82 (7%)	17	41

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	74	ARG
1	A	82	LEU
1	A	92	ARG
1	A	93	GLU
1	A	97	ARG
1	A	107	LEU
1	A	116	ARG
1	A	118	TYR
1	A	160	VAL
1	A	175	LEU
2	B	182	LEU
2	B	287	VAL
2	B	289	SER
2	B	326	ILE
2	B	337	LYS
2	B	342	GLN
2	B	347	LEU
2	B	348	ARG
2	B	355	LEU
2	B	387	LEU
1	C	10	ARG
1	C	21	LEU
1	C	41	SER
1	C	74	ARG
1	C	92	ARG
1	C	93	GLU
1	C	121	GLU
1	C	160	VAL
1	C	170	LYS
1	C	175	LEU
2	D	182	LEU
2	D	200	LEU
2	D	219	SER
2	D	226	LEU
2	D	266	ILE
2	D	294	THR
2	D	342	GLN
2	D	347	LEU
2	D	355	LEU
2	D	356	ARG
1	E	41	SER

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Mol	Chain	Res	Type
1	E	74	ARG
1	E	80	THR
1	E	89	ARG
1	E	92	ARG
1	E	108	ILE
1	E	113	VAL
1	E	152	ARG
1	E	160	VAL
1	E	175	LEU
3	F	200	LEU
3	F	214	ILE
3	F	216	THR
3	F	242	GLU
3	F	254	ASP
3	F	333	VAL
3	F	355	LEU
1	G	8	THR
1	G	39	THR
1	G	69	VAL
1	G	74	ARG
1	G	92	ARG
1	G	107	LEU
1	G	113	VAL
1	G	125	GLU
1	G	129	THR
1	G	150	ASP
1	G	160	VAL
1	G	170	LYS
1	G	175	LEU
2	H	182	LEU
2	H	200	LEU
2	H	216	THR
2	H	219	SER
2	H	221	SER
2	H	257	SER
2	H	275	ASP
2	H	317	THR
2	H	333	VAL
2	H	353	GLU
2	H	386	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

3 non-standard protein/DNA/RNA residues 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	TH5	B	181	2	9,9,10	1.86	2 (22%)	7,11,13	4.78	6 (85%)
2	TH5	D	181	2	9,9,10	1.69	2 (22%)	7,11,13	3.19	5 (71%)
2	TH5	H	181	2	9,9,10	2.66	2 (22%)	7,11,13	3.53	4 (57%)

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	TH5	B	181	2	-	0/9/10/12	0/0/0/0
2	TH5	D	181	2	-	0/9/10/12	0/0/0/0
2	TH5	H	181	2	-	0/9/10/12	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	181	TH5	OG1-CB	-6.06	1.36	1.46
2	B	181	TH5	O-C	-4.94	1.21	1.42
2	H	181	TH5	O-C	-4.77	1.21	1.42
2	D	181	TH5	O-C	-3.45	1.27	1.42
2	D	181	TH5	OG1-CB	-2.77	1.41	1.46
2	B	181	TH5	OG1-CB	-2.23	1.42	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	181	TH5	OG1-CB-CA	-9.41	90.93	107.27
2	H	181	TH5	OG1-CB-CA	-5.83	97.14	107.27
2	D	181	TH5	CB-OG1-CAH	-3.32	110.98	117.98
2	D	181	TH5	OG1-CB-CG2	-3.15	101.97	108.12
2	D	181	TH5	OG1-CB-CA	-2.92	102.20	107.27
2	H	181	TH5	OG1-CAH-OAD	-2.38	118.16	122.92
2	B	181	TH5	OAD-CAH-CAA	-2.11	117.03	124.85
2	D	181	TH5	O-C-CA	2.87	117.57	111.53
2	B	181	TH5	OG1-CB-CG2	3.39	114.75	108.12
2	B	181	TH5	CB-OG1-CAH	3.71	125.81	117.98
2	H	181	TH5	O-C-CA	3.90	119.73	111.53
2	B	181	TH5	OG1-CAH-CAA	4.45	119.50	111.10
2	B	181	TH5	O-C-CA	4.69	121.41	111.53
2	D	181	TH5	OG1-CAH-CAA	5.33	121.16	111.10
2	H	181	TH5	OG1-CAH-CAA	5.34	121.18	111.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	181	TH5	6	0
2	D	181	TH5	4	0
2	H	181	TH5	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	GLU	B	1601	-	4,8,9	1.50	1 (25%)	3,9,11	0.35	0
4	GLU	C	1602	-	4,8,9	1.44	0	3,9,11	0.55	0
5	ACT	E	1181	-	1,3,3	2.23	1 (100%)	0,3,3	0.00	-
5	ACT	G	1181	-	1,3,3	0.98	0	0,3,3	0.00	-

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	GLU	B	1601	-	-	0/3/7/9	0/0/0/0
4	GLU	C	1602	-	-	0/3/7/9	0/0/0/0
5	ACT	E	1181	-	-	0/0/0/0	0/0/0/0
5	ACT	G	1181	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	1181	ACT	CH3-C	2.23	1.51	1.48
4	B	1601	GLU	CB-CA	2.47	1.56	1.53

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1601	GLU	7	0
4	C	1602	GLU	5	0

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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	173/180 (96%)	-0.18	0 100 100	1, 9, 22, 29	0
1	C	173/180 (96%)	-0.10	1 (0%) 90 91	1, 7, 20, 24	0
1	E	173/180 (96%)	0.06	0 100 100	4, 15, 25, 28	0
1	G	173/180 (96%)	0.37	4 (2%) 64 64	6, 16, 29, 33	0
2	B	212/213 (99%)	-0.22	1 (0%) 91 93	1, 6, 19, 28	0
2	D	209/213 (98%)	-0.18	0 100 100	1, 7, 20, 29	0
2	H	194/213 (91%)	0.10	2 (1%) 84 85	6, 17, 32, 40	0
3	F	197/213 (92%)	-0.12	0 100 100	2, 14, 25, 31	0
All	All	1504/1572 (95%)	-0.04	8 (0%) 91 93	1, 11, 25, 40	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	336	PRO	4.6
1	G	22	ALA	3.3
1	G	134	ALA	2.9
2	H	195	ALA	2.8
1	C	134	ALA	2.6
1	G	88	ALA	2.5
1	G	135	GLY	2.1
2	B	342	GLN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	TH5	H	181	10/11	0.93	0.17	-	9,13,14,15	0
2	TH5	B	181	10/11	0.94	0.16	-	1,3,7,7	0
2	TH5	D	181	10/11	0.92	0.16	-	5,8,13,13	0

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
4	GLU	B	1601	9/10	0.82	0.26	4.04	13,15,19,21	0
4	GLU	C	1602	9/10	0.74	0.25	3.12	26,28,31,31	0
5	ACT	E	1181	4/4	0.94	0.16	0.65	16,17,17,17	0
5	ACT	G	1181	4/4	0.94	0.14	-1.29	15,16,16,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.