



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:54 PM UTC

PDB ID : 2EJW / pdb_00002ejw
Title : Homoserine Dehydrogenase from *Thermus thermophilus* HB8
Authors : Omi, R.; Goto, M.; Miyahara, I.; Hirotsu, K.
Deposited on : 2007-03-21
Resolution : 1.70 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

<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

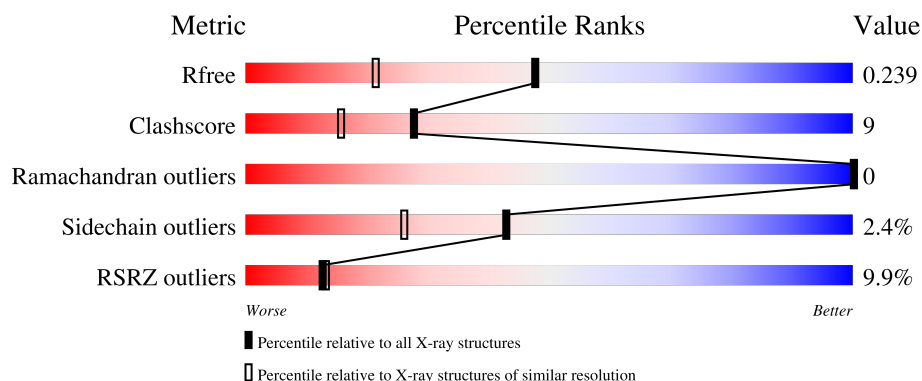
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 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	332	7% 81% 18% .
1	B	332	7% 80% 18% .
1	E	332	12% 81% 17% .
1	F	332	14% 77% 21% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10940 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.

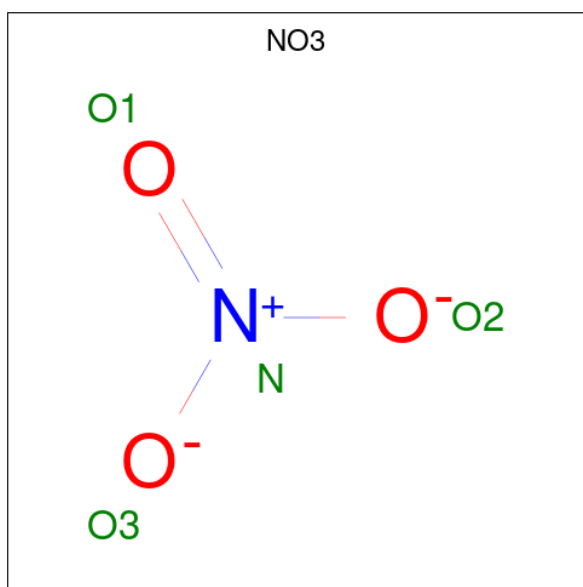
- Molecule 1 is a protein called Homoserine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2487	1594	438	451	4			
1	B	331	Total	C	N	O	S	0	0	0
			2493	1597	441	451	4			
1	E	331	Total	C	N	O	S	0	0	0
			2485	1593	438	450	4			
1	F	331	Total	C	N	O	S	0	0	0
			2487	1593	438	452	4			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

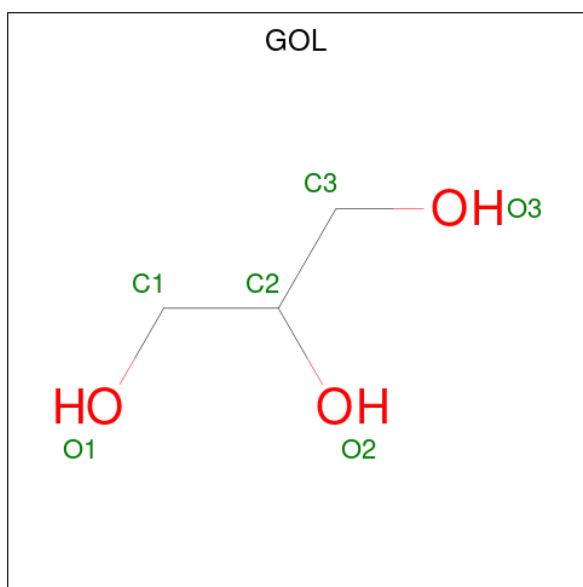
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		
2	B	2	Total	Mg	0	0
			2	2		
2	E	2	Total	Mg	0	0
			2	2		
2	F	1	Total	Mg	0	0
			1	1		

- Molecule 3 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	N	O	0	0
			4	1	3		
3	E	1	Total	N	O	0	0
			4	1	3		
3	F	1	Total	N	O	0	0
			4	1	3		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

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

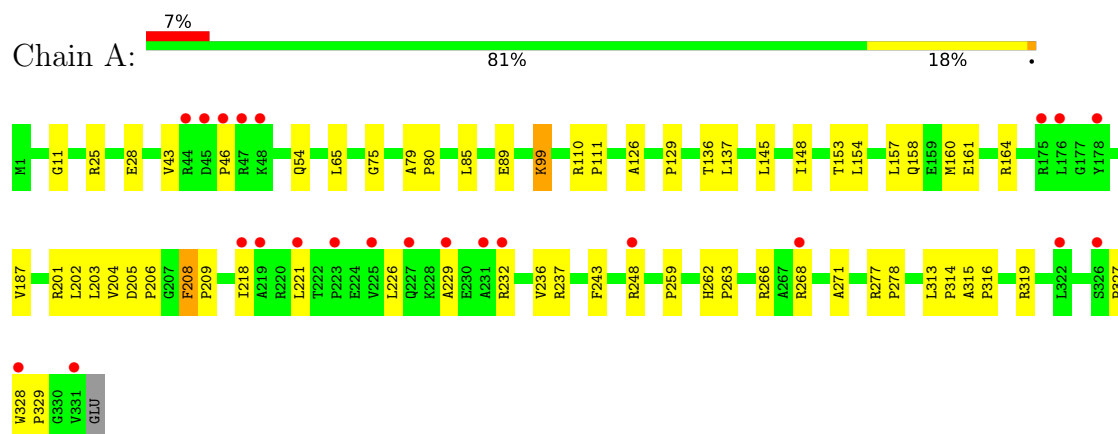
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	273	Total	O	0	0
			273	273		
5	B	234	Total	O	0	0
			234	234		
5	E	220	Total	O	0	0
			220	220		
5	F	230	Total	O	0	0
			230	230		

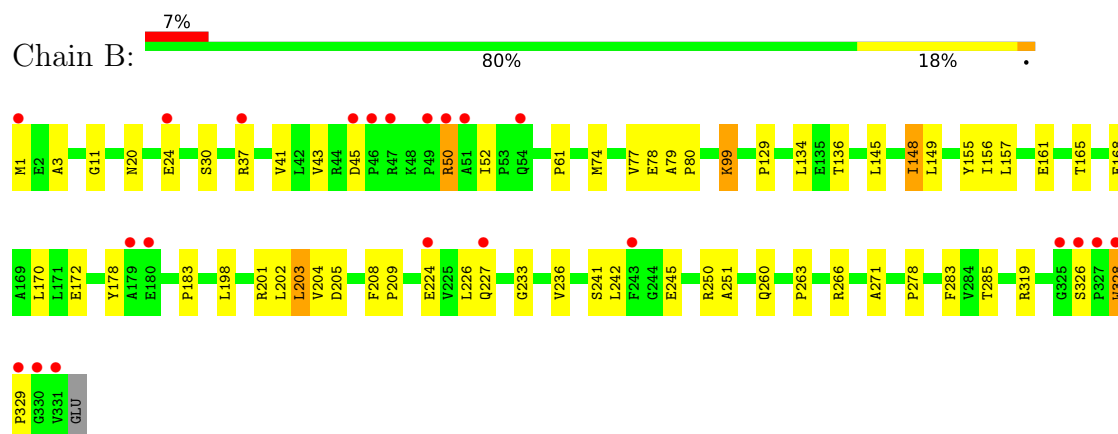
3 Residue-property plots [i](#)

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.

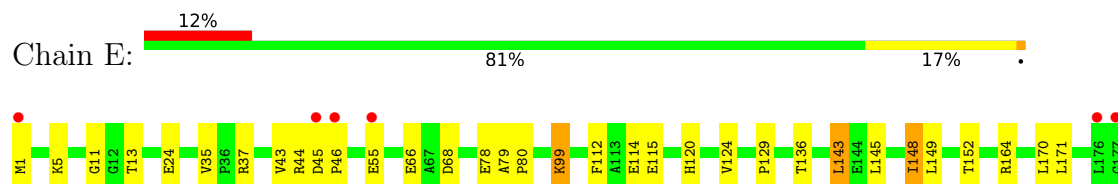
- Molecule 1: Homoserine dehydrogenase

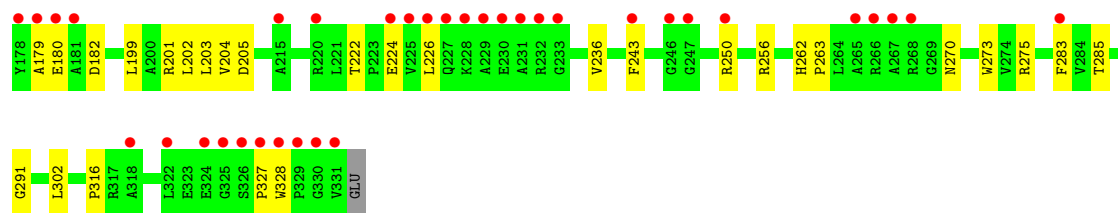


- Molecule 1: Homoserine dehydrogenase

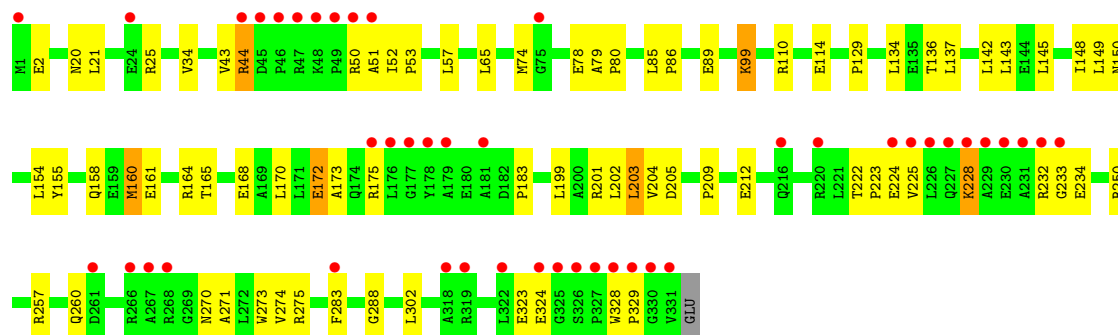
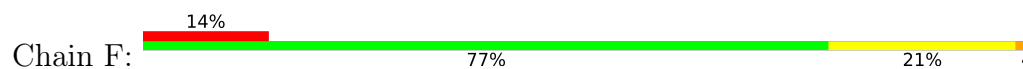


- Molecule 1: Homoserine dehydrogenase





● Molecule 1: Homoserine dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	165.44Å 173.60Å 97.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.36 – 1.70 41.36 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (41.36-1.70) 99.9 (41.36-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.72 (at 1.70Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.210 , 0.238 0.212 , 0.239	Depositor DCC
R_{free} test set	15266 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10940	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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: MG, GOL, NO3

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	0.40	0/2542	0.97	10/3462 (0.3%)
1	B	0.40	0/2548	0.93	8/3469 (0.2%)
1	E	0.39	0/2540	0.97	11/3459 (0.3%)
1	F	0.42	1/2542 (0.0%)	0.94	7/3461 (0.2%)
All	All	0.40	1/10172 (0.0%)	0.95	36/13851 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	160	MET	SD-CE	-5.97	1.64	1.79

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	204	VAL	N-CA-C	8.72	122.08	113.53
1	F	204	VAL	N-CA-C	8.45	121.81	113.53
1	E	204	VAL	N-CA-C	7.69	121.07	113.53
1	B	149	LEU	N-CA-C	7.65	122.61	113.28
1	B	77	VAL	N-CA-C	7.10	118.16	111.48
1	B	78	GLU	N-CA-C	6.87	118.42	111.07
1	E	148	ILE	N-CA-C	-6.85	95.82	106.72
1	F	65	LEU	N-CA-C	6.71	120.56	112.38
1	A	148	ILE	N-CA-C	-6.68	96.10	106.72
1	B	204	VAL	N-CA-C	6.65	120.05	113.53
1	A	65	LEU	N-CA-C	6.62	120.82	112.87
1	F	78	GLU	N-CA-C	6.53	118.95	111.11
1	E	182	ASP	CA-C-N	6.35	127.17	120.12
1	E	182	ASP	C-N-CA	6.35	127.17	120.12
1	E	78	GLU	N-CA-C	6.14	117.64	111.07
1	E	203	LEU	N-CA-C	6.04	120.72	112.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	203	LEU	N-CA-C	5.94	120.58	112.68
1	E	45	ASP	CA-C-N	5.85	125.72	119.28
1	E	45	ASP	C-N-CA	5.85	125.72	119.28
1	F	149	LEU	N-CA-C	5.84	120.10	112.92
1	B	203	LEU	N-CA-C	5.68	120.23	112.68
1	A	75	GLY	N-CA-C	-5.66	103.56	112.61
1	E	68	ASP	N-CA-C	-5.62	105.53	112.90
1	A	203	LEU	N-CA-C	5.58	120.08	112.04
1	B	148	ILE	N-CA-C	-5.54	97.81	109.34
1	B	50	ARG	N-CA-C	5.39	117.39	109.24
1	A	126	ALA	CB-CA-C	-5.34	110.40	116.54
1	E	120	HIS	N-CA-C	5.31	117.25	110.61
1	F	148	ILE	N-CA-C	-5.27	99.37	106.85
1	E	149	LEU	N-CA-C	5.26	119.96	113.50
1	B	77	VAL	CB-CA-C	-5.23	107.26	111.71
1	A	187	VAL	N-CA-C	5.19	119.75	113.00
1	F	143	LEU	N-CA-C	-5.18	105.79	112.68
1	A	271	ALA	N-CA-C	5.15	116.89	109.07
1	A	208	PHE	CA-C-N	5.02	125.17	119.90
1	A	208	PHE	C-N-CA	5.02	125.17	119.90

There are no chirality outliers.

There are no planarity outliers.

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	2487	0	2524	40	0
1	B	2493	0	2535	43	0
1	E	2485	0	2520	43	0
1	F	2487	0	2526	64	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
3	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	4	0	0	0	0
3	F	4	0	0	0	0
4	B	6	0	8	1	0
4	E	6	0	8	0	0
5	A	273	0	0	2	0
5	B	234	0	0	1	0
5	E	220	0	0	4	0
5	F	230	0	0	1	0
All	All	10940	0	10121	181	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 9.

All (181) 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:99:LYS:H	1:F:99:LYS:HD3	1.24	0.99
1:A:218:ILE:HG23	1:A:221:LEU:HD12	1.49	0.93
1:A:153:THR:HG22	1:A:218:ILE:HD11	1.60	0.84
1:B:99:LYS:H	1:B:99:LYS:HD3	1.41	0.84
1:A:46:PRO:HB2	1:A:54:GLN:HE22	1.45	0.81
1:F:20:ASN:HD21	1:F:51:ALA:HB1	1.47	0.78
1:B:20:ASN:O	1:B:24:GLU:HG3	1.85	0.77
1:B:3:ALA:HB2	1:B:37:ARG:NH1	2.01	0.76
1:F:225:VAL:HA	1:F:228:LYS:NZ	2.03	0.73
1:F:79:ALA:HB3	1:F:80:PRO:HD3	1.71	0.73
1:F:225:VAL:HA	1:F:228:LYS:HZ3	1.54	0.72
1:F:232:ARG:HB2	1:F:232:ARG:HH11	1.55	0.71
1:A:85:LEU:O	1:A:89:GLU:HG3	1.90	0.71
1:F:155:TYR:OH	1:F:172:GLU:HG2	1.91	0.71
1:A:79:ALA:HB3	1:A:80:PRO:HD3	1.75	0.69
1:B:157:LEU:O	1:B:161:GLU:HG3	1.94	0.68
1:E:79:ALA:HB3	1:E:80:PRO:HD3	1.77	0.66
1:E:226:LEU:CD1	1:E:236:VAL:HB	2.25	0.66
1:A:99:LYS:HD2	1:A:99:LYS:H	1.61	0.66
1:A:266:ARG:HD3	1:B:266:ARG:HD2	1.78	0.65
1:F:165:THR:OG1	1:F:168:GLU:HG3	1.96	0.64
1:A:248:ARG:HG3	1:A:248:ARG:HH11	1.60	0.64
1:F:99:LYS:HE3	5:F:2654:HOH:O	1.96	0.64
1:B:11:GLY:HA3	1:B:43:VAL:HG12	1.78	0.63
1:E:5:LYS:HE2	1:E:37:ARG:HH21	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:224:GLU:CD	1:E:224:GLU:H	2.08	0.62
1:F:232:ARG:HH12	1:F:234:GLU:CD	2.09	0.61
1:F:74:MET:HG3	1:F:80:PRO:HG3	1.83	0.61
1:A:154:LEU:O	1:A:158:GLN:HG3	2.00	0.61
1:E:201:ARG:HA	1:E:205:ASP:O	1.99	0.61
1:F:232:ARG:HH11	1:F:232:ARG:CB	2.14	0.61
1:F:234:GLU:OE1	1:F:257:ARG:HD3	2.01	0.60
1:E:164:ARG:HH11	1:E:164:ARG:HG3	1.66	0.60
1:F:201:ARG:HA	1:F:205:ASP:O	2.02	0.60
1:A:229:ALA:HA	1:A:232:ARG:HH21	1.65	0.60
1:A:157:LEU:O	1:A:161:GLU:HB2	2.02	0.60
1:A:277:ARG:HG3	1:A:277:ARG:HH11	1.66	0.60
1:F:20:ASN:ND2	1:F:51:ALA:HB1	2.17	0.60
1:F:228:LYS:NZ	1:F:228:LYS:HB3	2.19	0.57
1:E:222:THR:O	1:E:226:LEU:HD23	2.03	0.57
1:E:285:THR:HG21	5:E:2789:HOH:O	2.04	0.57
1:B:165:THR:OG1	1:B:168:GLU:HG3	2.05	0.57
1:E:263:PRO:HB3	1:E:283:PHE:CE2	2.39	0.57
1:A:25:ARG:NH2	1:A:28:GLU:OE1	2.37	0.56
1:B:79:ALA:HB3	1:B:80:PRO:HD3	1.87	0.56
1:B:3:ALA:HB1	1:B:37:ARG:HD2	1.87	0.56
1:E:273:TRP:CH2	1:E:275:ARG:HG3	2.41	0.56
1:F:172:GLU:OE1	1:F:175:ARG:NH2	2.38	0.55
1:F:85:LEU:O	1:F:89:GLU:HG3	2.07	0.55
1:F:228:LYS:HZ3	1:F:228:LYS:HB3	1.71	0.55
1:A:153:THR:O	1:A:157:LEU:HG	2.06	0.54
1:B:3:ALA:CB	1:B:37:ARG:HD2	2.37	0.54
1:B:226:LEU:CD2	1:B:236:VAL:HB	2.37	0.54
1:E:226:LEU:HD12	5:E:2780:HOH:O	2.06	0.54
1:F:232:ARG:CB	1:F:232:ARG:NH1	2.70	0.54
1:A:160:MET:HA	1:A:164:ARG:O	2.08	0.54
1:E:226:LEU:HD11	1:E:236:VAL:HB	1.88	0.54
1:F:223:PRO:HD2	1:F:224:GLU:OE2	2.08	0.53
1:B:148:ILE:O	1:B:148:ILE:HG13	2.08	0.53
1:A:328:TRP:HB3	1:A:329:PRO:HD2	1.91	0.53
1:B:156:ILE:HD11	1:B:170:LEU:HD13	1.90	0.52
1:E:143:LEU:HD23	1:E:275:ARG:NH2	2.25	0.52
1:F:222:THR:OG1	1:F:225:VAL:HG23	2.10	0.52
1:F:20:ASN:HD21	1:F:51:ALA:CB	2.20	0.52
1:E:226:LEU:HD13	1:E:236:VAL:HB	1.92	0.52
1:F:99:LYS:HD3	1:F:99:LYS:N	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:202:LEU:C	1:F:202:LEU:HD23	2.35	0.52
1:E:11:GLY:HA3	1:E:43:VAL:HG12	1.92	0.51
1:F:50:ARG:HG2	1:F:50:ARG:HH11	1.75	0.51
1:F:202:LEU:HD23	1:F:202:LEU:O	2.10	0.51
1:F:2:GLU:O	1:F:34:VAL:HA	2.11	0.51
1:F:85:LEU:HB2	1:F:86:PRO:HD3	1.93	0.51
1:F:228:LYS:O	1:F:232:ARG:HD3	2.11	0.51
1:F:43:VAL:HG11	1:F:57:LEU:HD22	1.93	0.50
1:A:218:ILE:HG22	1:A:218:ILE:O	2.11	0.50
1:A:243:PHE:CZ	1:A:327:PRO:HD3	2.46	0.50
1:F:150:ASN:C	1:F:270:ASN:HD21	2.19	0.50
1:E:112:PHE:HA	1:E:115:GLU:HG2	1.94	0.50
1:E:263:PRO:HB3	1:E:283:PHE:CZ	2.47	0.50
1:A:248:ARG:HG3	1:A:248:ARG:NH1	2.27	0.50
1:B:201:ARG:HA	1:B:205:ASP:O	2.11	0.49
1:E:285:THR:HG22	1:F:283:PHE:HB3	1.92	0.49
1:E:99:LYS:H	1:E:99:LYS:HD2	1.77	0.49
1:F:145:LEU:C	1:F:145:LEU:HD23	2.38	0.49
1:B:263:PRO:HB3	1:B:283:PHE:CE2	2.48	0.49
1:F:209:PRO:HG2	1:F:212:GLU:HG3	1.94	0.49
1:B:145:LEU:HB3	1:B:242:LEU:HB3	1.94	0.49
1:B:74:MET:HG3	1:B:80:PRO:HG3	1.95	0.48
1:A:99:LYS:H	1:A:99:LYS:CD	2.26	0.48
1:E:152:THR:OG1	1:E:179:ALA:HB2	2.12	0.48
1:B:1:MET:HE1	1:B:30:SER:HA	1.94	0.48
1:F:50:ARG:HG2	1:F:50:ARG:NH1	2.29	0.48
1:A:129:PRO:HG3	1:B:136:THR:HB	1.94	0.48
1:A:201:ARG:HA	1:A:205:ASP:O	2.13	0.48
1:A:319:ARG:O	1:A:319:ARG:HD2	2.13	0.48
1:B:271:ALA:HB2	1:B:285:THR:HG22	1.95	0.48
1:F:273:TRP:CH2	1:F:275:ARG:HG3	2.49	0.47
1:A:226:LEU:CD2	1:A:236:VAL:HB	2.45	0.47
1:E:164:ARG:HG3	1:E:164:ARG:NH1	2.29	0.47
1:B:3:ALA:HB2	1:B:37:ARG:HH11	1.74	0.47
1:E:136:THR:HB	1:F:129:PRO:HG3	1.97	0.47
1:F:160:MET:HE1	1:F:222:THR:HA	1.97	0.47
1:A:110:ARG:N	1:A:111:PRO:HD2	2.30	0.46
1:B:271:ALA:CB	1:B:285:THR:HG22	2.45	0.46
1:F:271:ALA:HB1	1:F:283:PHE:CE1	2.50	0.46
1:F:302:LEU:C	1:F:302:LEU:HD23	2.41	0.46
1:B:178:TYR:CE1	4:B:2501:GOL:H31	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:LEU:C	1:B:202:LEU:HD23	2.40	0.46
1:B:245:GLU:OE1	1:B:250:ARG:NH1	2.43	0.46
1:F:233:GLY:O	1:F:260:GLN:HG2	2.15	0.46
1:A:262:HIS:ND1	1:A:263:PRO:HD2	2.31	0.46
1:E:170:LEU:C	1:E:170:LEU:HD23	2.42	0.45
1:B:145:LEU:C	1:B:145:LEU:HD23	2.42	0.45
1:F:155:TYR:CD2	1:F:173:ALA:HB2	2.51	0.45
1:E:24:GLU:HA	1:E:24:GLU:OE2	2.16	0.45
1:E:148:ILE:O	1:E:270:ASN:HB3	2.16	0.45
1:B:266:ARG:HG3	5:B:2603:HOH:O	2.17	0.45
1:F:228:LYS:NZ	1:F:228:LYS:CB	2.80	0.45
1:E:302:LEU:C	1:E:302:LEU:HD23	2.42	0.45
1:F:209:PRO:HG2	1:F:212:GLU:CG	2.47	0.45
1:B:41:VAL:HG11	1:B:52:ILE:HD13	1.99	0.45
1:E:129:PRO:HG3	1:F:136:THR:HB	1.98	0.45
1:B:224:GLU:O	1:B:227:GLN:HG2	2.17	0.44
1:E:35:VAL:HG13	1:E:35:VAL:O	2.18	0.44
1:E:145:LEU:C	1:E:145:LEU:HD23	2.43	0.44
1:F:158:GLN:O	1:F:161:GLU:HG2	2.16	0.44
1:A:157:LEU:CD1	1:A:237:ARG:HA	2.46	0.44
1:E:13:THR:HG22	1:E:291:GLY:HA3	2.00	0.44
1:E:66:GLU:HB2	5:E:2675:HOH:O	2.18	0.44
1:A:99:LYS:HD2	1:A:99:LYS:N	2.28	0.44
1:B:61:PRO:HD3	1:E:171:LEU:HD12	2.00	0.44
1:B:134:LEU:HD22	1:B:203:LEU:HD22	2.00	0.43
1:E:256:ARG:HB3	1:E:328:TRP:HB2	2.00	0.43
1:F:224:GLU:H	1:F:224:GLU:CD	2.25	0.43
1:A:206:PRO:HD3	1:A:315:ALA:HB3	2.01	0.43
1:E:44:ARG:O	1:E:46:PRO:HD3	2.17	0.43
1:F:164:ARG:HH22	1:F:172:GLU:HB2	1.83	0.43
1:B:328:TRP:HD1	1:B:329:PRO:HD2	1.83	0.43
1:F:154:LEU:O	1:F:158:GLN:HG3	2.18	0.43
1:F:44:ARG:HH11	1:F:44:ARG:CG	2.32	0.43
1:E:55:GLU:CD	1:E:55:GLU:H	2.27	0.43
1:B:170:LEU:HD11	1:B:183:PRO:HG3	2.01	0.43
1:E:285:THR:CG2	1:F:283:PHE:HB3	2.48	0.43
1:E:224:GLU:CD	1:E:224:GLU:N	2.75	0.42
1:A:11:GLY:HA3	1:A:43:VAL:HG12	2.01	0.42
1:A:316:PRO:HD3	5:A:2747:HOH:O	2.18	0.42
1:A:136:THR:HB	1:B:129:PRO:HG3	2.01	0.42
1:A:208:PHE:HA	1:A:209:PRO:HD3	1.79	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:274:VAL:HG12	1:F:275:ARG:N	2.34	0.42
1:F:110:ARG:O	1:F:114:GLU:HG3	2.18	0.42
1:B:241:SER:O	1:B:251:ALA:HA	2.20	0.42
1:E:243:PHE:CE2	1:E:327:PRO:HB3	2.55	0.42
1:F:328:TRP:HB3	1:F:329:PRO:HD2	2.01	0.42
1:F:99:LYS:H	1:F:99:LYS:CD	2.06	0.42
1:B:233:GLY:HA2	1:B:260:GLN:NE2	2.35	0.42
1:B:263:PRO:HB3	1:B:283:PHE:CZ	2.55	0.42
1:E:124:VAL:HG11	1:E:199:LEU:HD21	2.02	0.42
1:E:202:LEU:C	1:E:202:LEU:HD23	2.45	0.42
1:E:316:PRO:HD3	5:E:2707:HOH:O	2.20	0.42
1:A:145:LEU:C	1:A:145:LEU:HD23	2.45	0.41
1:A:218:ILE:HG22	1:A:221:LEU:HB2	2.02	0.41
1:F:52:ILE:HA	1:F:53:PRO:HD3	1.95	0.41
1:F:170:LEU:HD11	1:F:183:PRO:HG3	2.01	0.41
1:B:155:TYR:OH	1:B:172:GLU:HG2	2.20	0.41
1:B:224:GLU:H	1:B:224:GLU:CD	2.28	0.41
1:B:328:TRP:CD1	1:B:329:PRO:HD2	2.55	0.41
1:A:266:ARG:HD3	1:B:266:ARG:CD	2.46	0.41
1:B:208:PHE:HA	1:B:209:PRO:HD3	1.88	0.41
1:E:243:PHE:CE1	1:E:250:ARG:HD3	2.55	0.41
1:F:270:ASN:OD1	1:F:288:GLY:HA3	2.21	0.41
1:B:45:ASP:O	1:B:50:ARG:NH1	2.53	0.41
1:A:277:ARG:HH11	1:A:277:ARG:CG	2.33	0.41
1:E:262:HIS:CD2	1:E:328:TRP:CH2	3.09	0.41
1:F:21:LEU:HD11	1:F:25:ARG:HH11	1.85	0.41
1:A:259:PRO:HG3	5:A:2777:HOH:O	2.20	0.40
1:F:142:LEU:HD22	1:F:203:LEU:HD23	2.01	0.40
1:F:160:MET:HE3	1:F:223:PRO:CD	2.51	0.40
1:F:134:LEU:CD1	1:F:199:LEU:HD22	2.51	0.40
1:A:202:LEU:C	1:A:202:LEU:HD23	2.47	0.40
1:A:313:LEU:HA	1:A:314:PRO:HD3	1.92	0.40
1:F:160:MET:HE3	1:F:223:PRO:HG3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	329/332 (99%)	321 (98%)	8 (2%)	0	100	100
1	B	329/332 (99%)	323 (98%)	6 (2%)	0	100	100
1	E	329/332 (99%)	322 (98%)	7 (2%)	0	100	100
1	F	329/332 (99%)	322 (98%)	7 (2%)	0	100	100
All	All	1316/1328 (99%)	1288 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	244/248 (98%)	240 (98%)	4 (2%)	55	41
1	B	245/248 (99%)	239 (98%)	6 (2%)	43	26
1	E	243/248 (98%)	238 (98%)	5 (2%)	47	30
1	F	245/248 (99%)	237 (97%)	8 (3%)	33	17
All	All	977/992 (98%)	954 (98%)	23 (2%)	43	26

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	LYS
1	A	137	LEU

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Mol	Chain	Res	Type
1	A	268	ARG
1	A	278	PRO
1	B	99	LYS
1	B	198	LEU
1	B	278	PRO
1	B	319	ARG
1	B	326	SER
1	B	328	TRP
1	E	1	MET
1	E	99	LYS
1	E	114	GLU
1	E	143	LEU
1	E	180	GLU
1	F	44	ARG
1	F	99	LYS
1	F	137	LEU
1	F	172	GLU
1	F	228	LYS
1	F	250	ARG
1	F	323	GLU
1	F	324	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	174	GLN
1	B	20	ASN
1	B	158	GLN
1	B	227	GLN
1	B	260	GLN
1	E	54	GLN
1	E	158	GLN
1	F	20	ASN
1	F	227	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 7 are monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GOL	E	2502	-	5,5,5	0.40	0	5,5,5	0.34	0
3	NO3	E	2601	-	1,3,3	0.19	0	0,3,3	-	-
4	GOL	B	2501	-	5,5,5	0.32	0	5,5,5	0.21	0
3	NO3	A	2602	-	1,3,3	0.34	0	0,3,3	-	-
3	NO3	F	2603	-	1,3,3	0.23	0	0,3,3	-	-

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	GOL	B	2501	-	-	2/4/4/4	-
4	GOL	E	2502	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	2501	GOL	C1-C2-C3-O3
4	B	2501	GOL	O2-C2-C3-O3
4	E	2502	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2501	GOL	1	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	331/332 (99%)	0.38	23 (6%) 23 25	9, 16, 32, 40	0
1	B	331/332 (99%)	0.43	22 (6%) 24 26	9, 16, 35, 45	0
1	E	331/332 (99%)	0.70	41 (12%) 8 8	10, 19, 37, 50	0
1	F	331/332 (99%)	0.64	45 (13%) 7 6	10, 17, 39, 46	0
All	All	1324/1328 (99%)	0.54	131 (9%) 13 13	9, 17, 36, 50	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	331	VAL	5.7
1	E	328	TRP	5.5
1	E	331	VAL	5.4
1	E	45	ASP	5.3
1	F	331	VAL	5.2
1	B	327	PRO	5.2
1	F	46	PRO	4.7
1	F	229	ALA	4.6
1	E	327	PRO	4.6
1	B	49	PRO	4.4
1	F	328	TRP	4.3
1	F	50	ARG	4.3
1	F	1	MET	4.3
1	F	268	ARG	4.3
1	A	331	VAL	4.3
1	E	329	PRO	4.0
1	E	176	LEU	3.9
1	B	328	TRP	3.9
1	F	231	ALA	3.9
1	E	322	LEU	3.9
1	F	176	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	E	179	ALA	3.8
1	F	179	ALA	3.7
1	A	44	ARG	3.6
1	F	178	TYR	3.6
1	F	181	ALA	3.6
1	F	226	LEU	3.6
1	B	329	PRO	3.5
1	E	247	GLY	3.5
1	A	218	ILE	3.5
1	E	330	GLY	3.5
1	F	319	ARG	3.5
1	A	46	PRO	3.5
1	F	75	GLY	3.4
1	E	226	LEU	3.4
1	F	327	PRO	3.4
1	A	328	TRP	3.4
1	F	44	ARG	3.4
1	A	223	PRO	3.3
1	F	45	ASP	3.3
1	A	45	ASP	3.3
1	F	330	GLY	3.3
1	B	180	GLU	3.2
1	A	231	ALA	3.2
1	B	326	SER	3.2
1	E	229	ALA	3.2
1	F	24	GLU	3.1
1	E	178	TYR	3.1
1	A	175	ARG	3.1
1	B	1	MET	3.0
1	F	228	LYS	3.0
1	F	49	PRO	3.0
1	E	233	GLY	3.0
1	F	47	ARG	3.0
1	E	318	ALA	3.0
1	A	47	ARG	3.0
1	E	227	GLN	2.9
1	F	48	LYS	2.9
1	E	268	ARG	2.9
1	B	37	ARG	2.9
1	B	51	ALA	2.9
1	F	177	GLY	2.9
1	B	50	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	179	ALA	2.8
1	E	326	SER	2.8
1	B	47	ARG	2.8
1	E	1	MET	2.8
1	E	267	ALA	2.8
1	F	233	GLY	2.8
1	B	330	GLY	2.8
1	E	246	GLY	2.8
1	E	325	GLY	2.8
1	F	325	GLY	2.8
1	F	232	ARG	2.8
1	A	248	ARG	2.7
1	B	46	PRO	2.7
1	F	322	LEU	2.7
1	E	250	ARG	2.7
1	E	224	GLU	2.7
1	A	178	TYR	2.7
1	E	230	GLU	2.6
1	F	329	PRO	2.6
1	E	324	GLU	2.6
1	E	181	ALA	2.6
1	E	180	GLU	2.6
1	A	268	ARG	2.6
1	E	232	ARG	2.6
1	F	318	ALA	2.6
1	B	224	GLU	2.6
1	B	45	ASP	2.5
1	A	48	LYS	2.5
1	E	46	PRO	2.5
1	A	176	LEU	2.5
1	F	326	SER	2.5
1	E	243	PHE	2.5
1	A	225	VAL	2.5
1	F	225	VAL	2.5
1	E	228	LYS	2.4
1	B	227	GLN	2.4
1	F	227	GLN	2.4
1	E	266	ARG	2.4
1	F	175	ARG	2.4
1	E	231	ALA	2.3
1	F	220	ARG	2.3
1	A	229	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	224	GLU	2.3
1	F	261	ASP	2.3
1	F	267	ALA	2.3
1	A	232	ARG	2.3
1	B	325	GLY	2.2
1	E	283	PHE	2.2
1	F	266	ARG	2.2
1	A	322	LEU	2.2
1	B	243	PHE	2.2
1	E	225	VAL	2.2
1	F	51	ALA	2.2
1	A	227	GLN	2.2
1	A	221	LEU	2.1
1	F	283	PHE	2.1
1	F	324	GLU	2.1
1	E	55	GLU	2.1
1	F	230	GLU	2.1
1	A	219	ALA	2.1
1	B	24	GLU	2.1
1	A	326	SER	2.1
1	E	215	ALA	2.0
1	F	216	GLN	2.0
1	B	54	GLN	2.0
1	E	177	GLY	2.0
1	E	220	ARG	2.0
1	E	265	ALA	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($\text{\AA}^2$)	Q<0.9
3	NO3	E	2601	4/4	0.82	0.18	23,26,29,30	0
4	GOL	E	2502	6/6	0.86	0.14	28,32,33,34	0
3	NO3	F	2603	4/4	0.88	0.14	23,26,27,28	0
4	GOL	B	2501	6/6	0.89	0.12	24,27,28,31	0
3	NO3	A	2602	4/4	0.91	0.11	18,19,21,22	0
2	MG	E	2007	1/1	0.94	0.07	28,28,28,28	0
2	MG	A	2002	1/1	0.96	0.05	20,20,20,20	0
2	MG	F	2005	1/1	0.98	0.04	11,11,11,11	0
2	MG	B	2004	1/1	0.98	0.05	11,11,11,11	0
2	MG	B	2001	1/1	0.98	0.04	15,15,15,15	0
2	MG	E	2006	1/1	0.99	0.05	11,11,11,11	0
2	MG	A	2003	1/1	0.99	0.05	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.