



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:56 AM UTC

PDB ID : 3DLG / pdb_00003dlg
Title : Crystal structure of hiv-1 reverse transcriptase in complex with GW564511.
Authors : Ren, J.; Chamberlain, P.P.; Stammers, D.K.
Deposited on : 2008-06-27
Resolution : 2.20 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

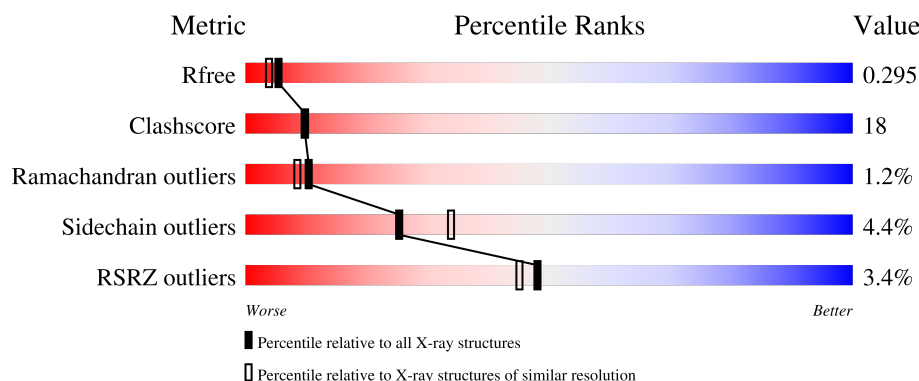
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	560	
2	B	440	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7928 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 REVERSE TRANSCRIPTASE/RIBONUCLEASE H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	532	Total	C	N	O	S	0	0	0
			4355	2822	723	802	8			

- Molecule 2 is a protein called P51 RT.

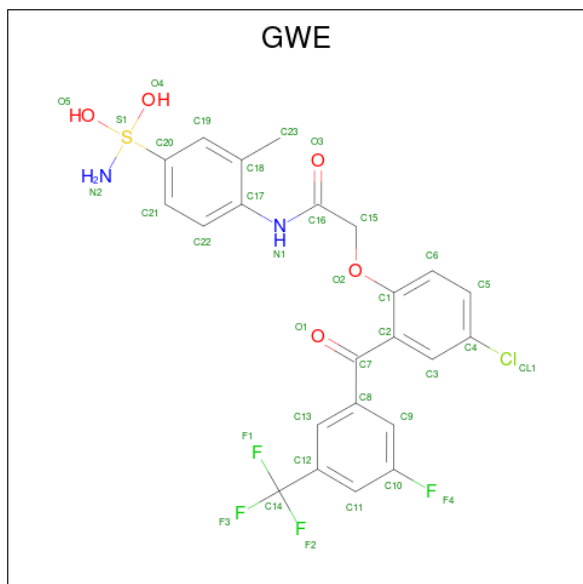
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	400	Total	C	N	O	S	0	0	0
			3303	2148	550	598	7			

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is N-{4-[amino(dihydroxy)-lambda 4 -sulfanyl]-2-methylphenyl}-2-(4-chloro-2-{[3-fluoro-5-(trifluoromethyl)phenyl]carbonyl}phenoxy)acetamide (CCD ID: GWE) (formula: $C_{23}H_{19}ClF_4N_2O_5S$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	Cl	F	N	O	S		
4	A	1	36	23	1	4	2	5	1	0	0

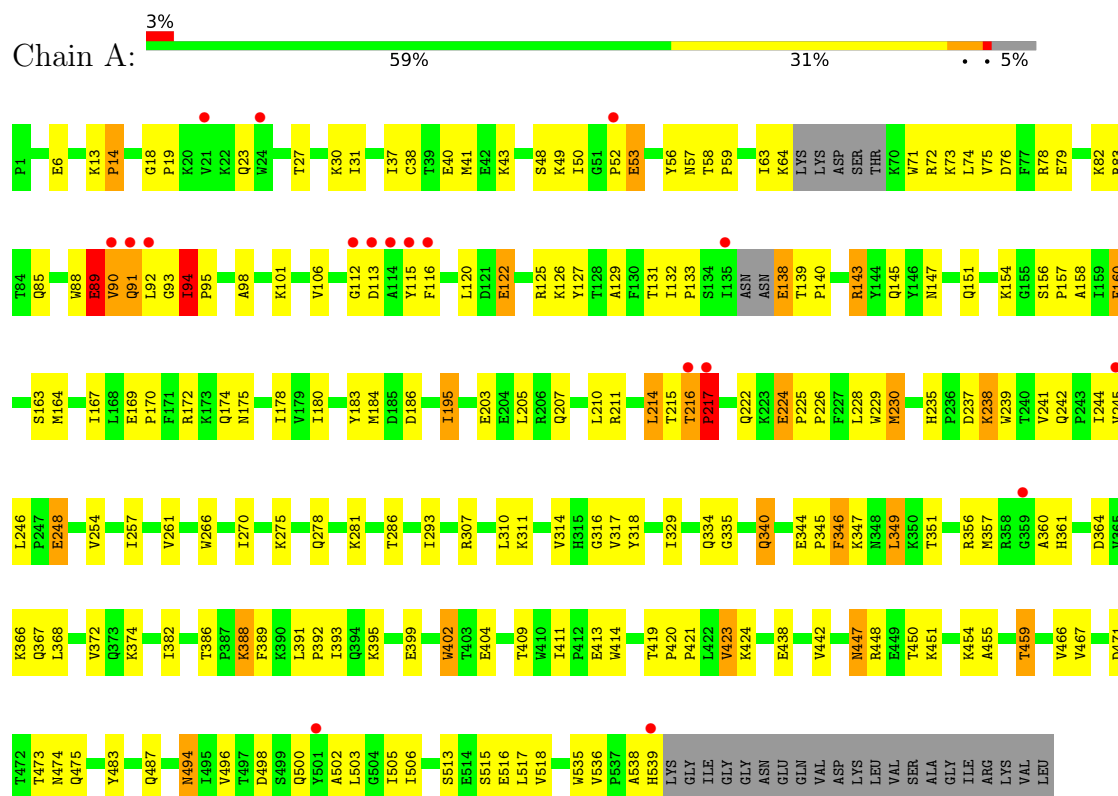
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	140	Total	O	0	0
			140	140		
5	B	84	Total	O	0	0
			84	84		

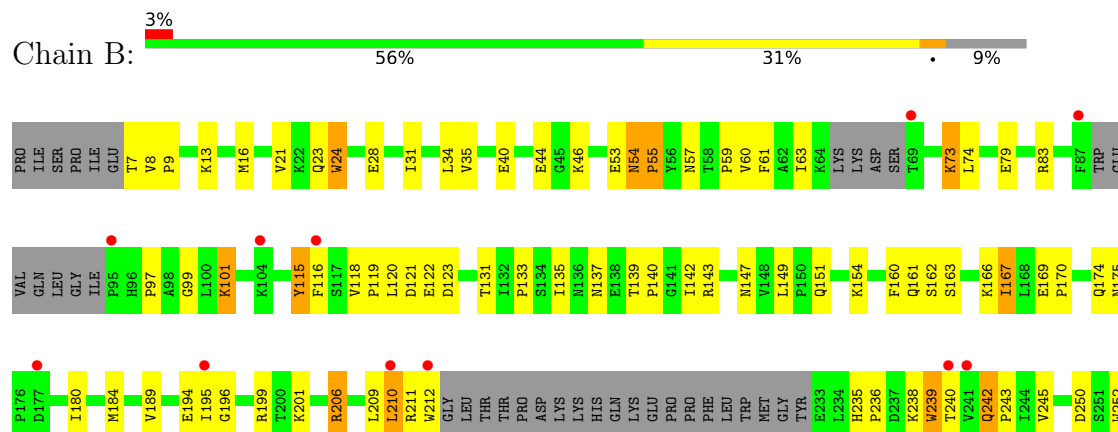
3 Residue-property plots

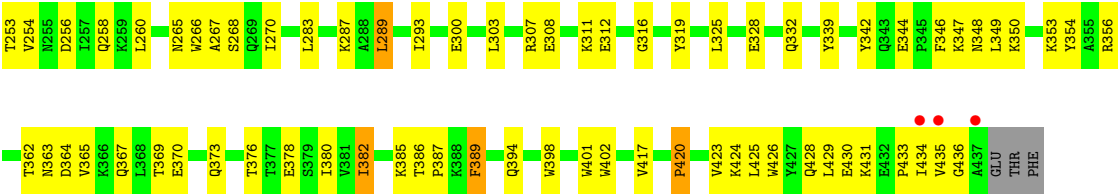
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: REVERSE TRANSCRIPTASE/RIBONUCLEASE H



• Molecule 2: P51 RT





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	137.22Å 109.52Å 72.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.43 – 2.20 29.43 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.43-2.20) 97.9 (29.43-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.299 0.217 , 0.295	Depositor DCC
R_{free} test set	2769 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	46.8	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7928	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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: PO4, CSD, GWE

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	0.47	0/4462	0.98	24/6064 (0.4%)
2	B	0.43	0/3394	0.96	20/4608 (0.4%)
All	All	0.45	0/7856	0.97	44/10672 (0.4%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	THR	CA-C-N	8.50	130.46	119.84
1	A	216	THR	C-N-CA	8.50	130.46	119.84
1	A	224	GLU	CA-C-N	8.00	125.50	119.66
1	A	224	GLU	C-N-CA	8.00	125.50	119.66
2	B	115	TYR	N-CA-C	7.95	120.02	111.36
2	B	402	TRP	N-CA-C	7.06	118.97	111.28
2	B	401	TRP	N-CA-C	6.41	123.03	114.12
2	B	349	LEU	N-CA-C	-6.33	104.61	112.90
2	B	417	VAL	N-CA-C	6.24	117.69	108.65
1	A	245	VAL	N-CA-C	6.21	117.05	107.99
1	A	349	LEU	N-CA-C	-6.21	103.82	112.45
2	B	382	ILE	N-CA-C	6.12	117.67	111.00
1	A	494	ASN	N-CA-C	-6.00	99.92	109.59
1	A	13	LYS	CA-C-N	5.98	127.32	119.84
1	A	13	LYS	C-N-CA	5.98	127.32	119.84
1	A	388	LYS	N-CA-C	-5.96	99.53	109.24
2	B	147	ASN	N-CA-C	-5.93	106.59	113.88
2	B	118	VAL	N-CA-C	5.86	115.35	108.96
1	A	154	LYS	N-CA-C	5.85	118.13	111.11
1	A	49	LYS	N-CA-C	-5.69	101.43	109.96
2	B	312	GLU	N-CA-C	5.61	114.76	108.25
2	B	250	ASP	N-CA-C	-5.61	104.18	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	THR	N-CA-C	5.60	118.22	108.76
1	A	516	GLU	N-CA-C	-5.59	105.27	111.36
2	B	428	GLN	N-CA-C	-5.55	105.77	112.54
1	A	217	PRO	N-CA-C	5.54	123.88	112.47
2	B	363	ASN	N-CA-C	5.51	119.39	107.49
1	A	186	ASP	N-CA-C	5.50	118.69	109.95
2	B	242	GLN	CA-C-N	5.50	125.69	120.31
2	B	242	GLN	C-N-CA	5.50	125.69	120.31
2	B	420	PRO	N-CA-C	-5.39	104.12	110.70
2	B	316	GLY	N-CA-C	5.36	123.16	114.90
1	A	389	PHE	N-CA-C	5.34	118.94	109.96
2	B	389	PHE	N-CA-C	5.32	118.73	110.17
1	A	447	ASN	N-CA-C	-5.24	101.23	109.25
1	A	382	ILE	N-CA-C	5.21	115.42	110.53
1	A	94	ILE	CA-C-N	5.12	125.04	119.76
1	A	94	ILE	C-N-CA	5.12	125.04	119.76
1	A	195	ILE	N-CA-C	5.11	116.44	110.62
2	B	54	ASN	CA-C-N	5.10	125.93	120.47
2	B	54	ASN	C-N-CA	5.10	125.93	120.47
2	B	131	THR	N-CA-C	5.06	117.31	108.76
1	A	92	LEU	N-CA-C	5.05	117.90	111.69
1	A	364	ASP	N-CA-C	5.04	116.78	111.28

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	4355	0	4398	164	0
2	B	3303	0	3338	118	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	36	0	17	3	0
5	A	140	0	0	3	0
5	B	84	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7928	0	7753	272	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 18.

All (272) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:353:LYS:HE3	2:B:430:GLU:HB3	1.46	0.97
2:B:242:GLN:HE21	2:B:243:PRO:HD2	1.28	0.96
1:A:270:ILE:HG22	1:A:314:VAL:HG11	1.48	0.93
1:A:216:THR:OG1	1:A:217:PRO:HD2	1.73	0.89
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.56	0.88
2:B:175:ASN:HD21	2:B:201:LYS:NZ	1.73	0.86
1:A:448:ARG:NE	1:A:474:ASN:H	1.74	0.86
2:B:195:ILE:HB	2:B:199:ARG:HH12	1.41	0.83
2:B:242:GLN:NE2	2:B:243:PRO:HD2	1.95	0.80
1:A:106:VAL:H	4:A:999:GWE:HNS	1.32	0.77
2:B:235:HIS:O	2:B:238:LYS:HG2	1.86	0.76
2:B:161:GLN:HE21	2:B:161:GLN:HA	1.50	0.75
1:A:448:ARG:HH21	1:A:475:GLN:H	1.34	0.75
2:B:175:ASN:HD21	2:B:201:LYS:HZ2	1.34	0.74
1:A:448:ARG:NH2	1:A:475:GLN:H	1.86	0.73
1:A:115:TYR:HB3	1:A:215:THR:O	1.90	0.72
2:B:161:GLN:HA	2:B:161:GLN:NE2	2.04	0.72
2:B:13:LYS:HB2	2:B:16:MET:HG3	1.70	0.71
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.24	0.71
2:B:356:ARG:HB2	2:B:367:GLN:HG2	1.71	0.71
1:A:126:LYS:HE2	1:A:127:TYR:CZ	2.24	0.71
1:A:239:TRP:CZ2	1:A:316:GLY:HA3	2.25	0.71
2:B:235:HIS:HB2	2:B:238:LYS:HE3	1.71	0.70
1:A:160:PHE:C	1:A:160:PHE:CD2	2.70	0.70
1:A:122:GLU:H	1:A:122:GLU:CD	1.98	0.69
2:B:63:ILE:HD13	2:B:74:LEU:HD22	1.73	0.69
1:A:393:ILE:HB	1:A:423:VAL:HG13	1.72	0.69
2:B:79:GLU:O	2:B:83:ARG:HG3	1.93	0.68
1:A:447:ASN:HB3	1:A:450:THR:OG1	1.92	0.68
1:A:448:ARG:HG3	1:A:448:ARG:HH11	1.59	0.68
1:A:50:ILE:HG21	1:A:145:GLN:HG2	1.75	0.67
2:B:376:THR:CG2	2:B:386:THR:HG22	2.24	0.67
2:B:7:THR:HG22	2:B:119:PRO:HG2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:LYS:HB2	1:A:316:GLY:O	1.95	0.67
1:A:183:TYR:HE1	1:A:184:MET:HE2	1.59	0.66
1:A:30:LYS:HE2	1:A:71:TRP:CH2	2.31	0.65
1:A:41:MET:HE1	1:A:73:LYS:HE2	1.79	0.65
2:B:195:ILE:HB	2:B:199:ARG:NH1	2.12	0.65
1:A:448:ARG:CZ	1:A:474:ASN:H	2.09	0.64
1:A:515:SER:HB3	1:A:518:VAL:HG23	1.78	0.64
1:A:448:ARG:HE	1:A:474:ASN:H	1.42	0.64
1:A:340:GLN:HB3	1:A:351:THR:HG22	1.80	0.64
2:B:116:PHE:HZ	2:B:151:GLN:NE2	1.96	0.64
2:B:151:GLN:HG3	5:B:1085:HOH:O	1.97	0.63
1:A:94:ILE:HD12	1:A:230:MET:HE2	1.80	0.63
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.81	0.63
1:A:211:ARG:HG2	1:A:211:ARG:HH11	1.64	0.63
2:B:376:THR:O	2:B:380:ILE:HG13	1.98	0.63
1:A:239:TRP:CE2	1:A:316:GLY:HA3	2.35	0.62
1:A:424:LYS:HB2	5:A:1067:HOH:O	1.98	0.62
1:A:503:LEU:HD13	1:A:535:TRP:CG	2.34	0.62
1:A:57:ASN:HD22	1:A:143:ARG:HH12	1.47	0.62
1:A:195:ILE:HG23	5:A:1023:HOH:O	1.99	0.62
2:B:350:LYS:HE2	2:B:378:GLU:OE1	1.99	0.62
2:B:73:LYS:NZ	2:B:73:LYS:HB3	2.15	0.62
1:A:160:PHE:C	1:A:160:PHE:HD2	2.07	0.62
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.82	0.62
1:A:203:GLU:HG3	1:A:207:GLN:HE21	1.65	0.61
1:A:257:ILE:O	1:A:261:VAL:HG23	2.00	0.61
2:B:376:THR:HG23	2:B:386:THR:HG22	1.81	0.61
2:B:369:THR:O	2:B:373:GLN:HG3	1.99	0.61
1:A:329:ILE:O	1:A:392:PRO:HD3	2.01	0.61
1:A:270:ILE:CG2	1:A:314:VAL:HG11	2.27	0.61
1:A:126:LYS:HE2	1:A:127:TYR:OH	2.01	0.60
1:A:160:PHE:HD2	1:A:160:PHE:O	1.84	0.60
1:A:366:LYS:NZ	2:B:394:GLN:NE2	2.49	0.60
1:A:335:GLY:HA3	1:A:356:ARG:CD	2.32	0.60
2:B:303:LEU:O	2:B:307:ARG:HG3	2.02	0.59
1:A:90:VAL:O	1:A:91:GLN:C	2.46	0.59
2:B:344:GLU:HB2	2:B:347:LYS:HD2	1.83	0.59
1:A:120:LEU:HD23	1:A:125:ARG:HG2	1.84	0.59
1:A:344:GLU:HG3	1:A:347:LYS:HD2	1.86	0.58
1:A:143:ARG:HH11	1:A:143:ARG:HG2	1.67	0.58
1:A:37:ILE:HG22	1:A:41:MET:HE3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:169:GLU:HB2	2:B:170:PRO:HD3	1.85	0.58
1:A:64:LYS:HA	1:A:72:ARG:HG3	1.86	0.57
1:A:448:ARG:HE	1:A:473:THR:HB	1.68	0.57
1:A:346:PHE:N	1:A:346:PHE:CD2	2.71	0.57
2:B:209:LEU:C	2:B:211:ARG:H	2.12	0.57
1:A:225:PRO:HA	1:A:226:PRO:C	2.29	0.56
2:B:163:SER:O	2:B:167:ILE:HG12	2.05	0.56
1:A:76:ASP:OD2	1:A:78:ARG:HG3	2.05	0.56
2:B:160:PHE:O	2:B:160:PHE:CD1	2.59	0.56
1:A:74:LEU:HD23	1:A:75:VAL:N	2.20	0.56
2:B:206:ARG:O	2:B:210:LEU:N	2.35	0.56
2:B:245:VAL:HG23	2:B:431:LYS:HB2	1.89	0.55
1:A:366:LYS:HZ2	2:B:394:GLN:NE2	2.03	0.55
2:B:194:GLU:CD	2:B:196:GLY:H	2.13	0.55
1:A:183:TYR:CE1	1:A:184:MET:HE2	2.41	0.55
2:B:101:LYS:O	2:B:236:PRO:HB2	2.07	0.55
2:B:13:LYS:HB2	2:B:16:MET:CG	2.36	0.55
1:A:254:VAL:HG22	1:A:286:THR:HG21	1.88	0.54
1:A:116:PHE:CE2	1:A:214:LEU:HD11	2.42	0.54
1:A:360:ALA:O	1:A:513:SER:HB2	2.08	0.54
1:A:388:LYS:HE2	1:A:413:GLU:OE1	2.07	0.54
2:B:139:THR:HG23	2:B:140:PRO:HD2	1.89	0.54
1:A:90:VAL:O	1:A:90:VAL:HG23	2.08	0.53
2:B:319:TYR:OH	2:B:385:LYS:HD3	2.07	0.53
2:B:167:ILE:HG22	2:B:212:TRP:CZ2	2.44	0.53
1:A:217:PRO:HB2	1:A:222:GLN:HE21	1.73	0.53
2:B:31:ILE:O	2:B:35:VAL:HG23	2.08	0.53
1:A:409:THR:O	2:B:364:ASP:HB2	2.09	0.53
1:A:183:TYR:CD1	1:A:184:MET:HG3	2.44	0.52
1:A:37:ILE:CG2	1:A:41:MET:HE3	2.39	0.52
1:A:356:ARG:HG2	1:A:356:ARG:HH11	1.74	0.52
1:A:395:LYS:HD3	1:A:414:TRP:CZ2	2.45	0.52
1:A:502:ALA:HA	1:A:505:ILE:HD12	1.92	0.52
2:B:175:ASN:ND2	2:B:201:LYS:NZ	2.53	0.52
1:A:335:GLY:HA3	1:A:356:ARG:HD3	1.92	0.52
1:A:40:GLU:O	1:A:43:LYS:HG2	2.10	0.52
1:A:366:LYS:NZ	2:B:394:GLN:HE21	2.07	0.51
2:B:252:TRP:CZ3	2:B:260:LEU:HD22	2.45	0.51
1:A:515:SER:HB3	1:A:518:VAL:CG2	2.39	0.51
2:B:53:GLU:O	2:B:55:PRO:HD3	2.10	0.51
2:B:344:GLU:CB	2:B:347:LYS:HD2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:ASP:HB3	1:A:151:GLN:NE2	2.26	0.51
1:A:246:LEU:O	1:A:307:ARG:NH1	2.35	0.51
1:A:18:GLY:HA3	1:A:56:TYR:CD1	2.46	0.51
2:B:57:ASN:HD22	2:B:143:ARG:NH1	2.09	0.51
1:A:50:ILE:CG2	1:A:145:GLN:HG2	2.40	0.51
1:A:79:GLU:O	1:A:83:ARG:HG3	2.10	0.51
1:A:58:THR:HG23	1:A:59:PRO:HD2	1.92	0.51
1:A:58:THR:HG23	1:A:76:ASP:O	2.11	0.51
2:B:160:PHE:O	2:B:160:PHE:HD1	1.94	0.51
1:A:94:ILE:HD11	1:A:230:MET:HG3	1.94	0.50
1:A:125:ARG:HD3	1:A:147:ASN:HA	1.92	0.50
1:A:402:TRP:CZ2	2:B:362:THR:HA	2.47	0.50
2:B:139:THR:HG22	2:B:140:PRO:O	2.10	0.50
1:A:307:ARG:O	1:A:311:LYS:HD3	2.10	0.50
1:A:237:ASP:OD2	1:A:237:ASP:N	2.44	0.50
1:A:448:ARG:HE	1:A:474:ASN:N	2.09	0.50
2:B:425:LEU:O	2:B:429:LEU:HD13	2.12	0.50
2:B:354:TYR:OH	2:B:370:GLU:HB3	2.10	0.50
2:B:239:TRP:CH2	2:B:378:GLU:HG2	2.46	0.50
1:A:293:ILE:N	1:A:293:ILE:HD12	2.27	0.50
2:B:24:TRP:HH2	2:B:61:PHE:CD2	2.29	0.50
2:B:369:THR:HG22	2:B:398:TRP:CZ3	2.47	0.50
2:B:162:SER:O	2:B:166:LYS:HG3	2.12	0.50
2:B:24:TRP:CH2	2:B:61:PHE:CD2	3.00	0.49
1:A:244:ILE:HB	1:A:310:LEU:HD13	1.94	0.49
2:B:154:LYS:HA	2:B:184:MET:HE1	1.92	0.49
1:A:494:ASN:HB3	2:B:289:LEU:HD22	1.94	0.49
2:B:34:LEU:CD2	2:B:73:LYS:HG3	2.42	0.49
1:A:217:PRO:HB2	1:A:222:GLN:NE2	2.27	0.49
1:A:27:THR:O	1:A:31:ILE:HG13	2.13	0.49
2:B:206:ARG:O	2:B:210:LEU:HB2	2.12	0.49
1:A:420:PRO:HA	1:A:421:PRO:C	2.37	0.49
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.52	0.49
1:A:346:PHE:N	1:A:346:PHE:HD2	2.10	0.49
1:A:372:VAL:HG11	1:A:411:ILE:HG23	1.94	0.49
1:A:335:GLY:HA3	1:A:356:ARG:HD2	1.95	0.48
2:B:115:TYR:HB3	2:B:149:LEU:CB	2.33	0.48
1:A:183:TYR:CE1	1:A:184:MET:HG3	2.48	0.48
1:A:138:GLU:OE1	1:A:139:THR:HG23	2.13	0.48
1:A:139:THR:HB	1:A:140:PRO:CD	2.43	0.48
2:B:254:VAL:O	2:B:258:GLN:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:387:PRO:HG2	2:B:389:PHE:CE1	2.48	0.48
2:B:434:ILE:HG22	2:B:435:VAL:HG13	1.96	0.48
1:A:63:ILE:HD12	1:A:64:LYS:N	2.28	0.48
1:A:442:VAL:HG22	1:A:496:VAL:O	2.14	0.48
1:A:211:ARG:HG2	1:A:211:ARG:NH1	2.25	0.48
2:B:8:VAL:HG23	2:B:9:PRO:HD2	1.94	0.48
2:B:206:ARG:NH1	2:B:206:ARG:HG2	2.28	0.48
1:A:335:GLY:O	1:A:356:ARG:HA	2.13	0.48
1:A:448:ARG:HE	1:A:473:THR:CA	2.27	0.48
1:A:454:LYS:HA	1:A:467:VAL:O	2.15	0.47
2:B:265:ASN:O	2:B:268:SER:HB3	2.14	0.47
2:B:235:HIS:CB	2:B:238:LYS:HE3	2.40	0.47
2:B:426:TRP:O	2:B:429:LEU:HB2	2.14	0.47
1:A:139:THR:HB	1:A:140:PRO:HD2	1.96	0.47
1:A:57:ASN:HB2	1:A:143:ARG:HH22	1.80	0.47
2:B:238:LYS:O	2:B:240:THR:N	2.48	0.47
1:A:170:PRO:O	1:A:174:GLN:HG3	2.15	0.47
1:A:172:ARG:NH2	1:A:180:ILE:O	2.48	0.46
2:B:120:LEU:O	2:B:121:ASP:C	2.56	0.46
1:A:535:TRP:HH2	5:B:1183:HOH:O	1.97	0.46
2:B:73:LYS:HB3	2:B:73:LYS:HZ3	1.80	0.46
2:B:175:ASN:ND2	2:B:201:LYS:HZ2	2.07	0.46
1:A:356:ARG:HG2	1:A:356:ARG:NH1	2.29	0.46
2:B:54:ASN:O	2:B:143:ARG:NH2	2.48	0.46
1:A:216:THR:O	1:A:217:PRO:O	2.34	0.46
2:B:209:LEU:C	2:B:211:ARG:N	2.74	0.46
2:B:61:PHE:CZ	2:B:74:LEU:HD23	2.51	0.46
2:B:180:ILE:HG12	2:B:189:VAL:HG13	1.98	0.46
1:A:183:TYR:CD1	1:A:229:TRP:HH2	2.34	0.46
1:A:334:GLN:O	1:A:356:ARG:HD3	2.15	0.46
2:B:211:ARG:HG3	2:B:212:TRP:CD1	2.51	0.45
1:A:329:ILE:HD12	1:A:391:LEU:CD2	2.46	0.45
1:A:399:GLU:OE2	1:A:402:TRP:HZ3	1.98	0.45
2:B:175:ASN:HD21	2:B:201:LYS:HZ1	1.58	0.45
2:B:433:PRO:HB3	2:B:436:GLY:O	2.16	0.45
1:A:23:GLN:OE1	1:A:59:PRO:HA	2.17	0.45
2:B:287:LYS:CD	2:B:293:ILE:HD11	2.47	0.45
2:B:308:GLU:O	2:B:311:LYS:HB2	2.16	0.45
1:A:402:TRP:CH2	2:B:362:THR:HA	2.52	0.45
2:B:308:GLU:OE1	2:B:311:LYS:HG3	2.16	0.45
1:A:74:LEU:HD23	1:A:74:LEU:C	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.97	0.45
1:A:244:ILE:HD11	1:A:266:TRP:HZ3	1.81	0.45
2:B:170:PRO:O	2:B:174:GLN:HG3	2.17	0.44
2:B:245:VAL:CG2	2:B:431:LYS:HB2	2.47	0.44
1:A:402:TRP:CD1	1:A:402:TRP:C	2.94	0.44
1:A:419:THR:HG22	5:A:1134:HOH:O	2.16	0.44
1:A:90:VAL:CG1	1:A:158:ALA:HB2	2.48	0.44
2:B:122:GLU:HG3	2:B:123:ASP:N	2.32	0.44
2:B:206:ARG:HG2	2:B:206:ARG:HH11	1.81	0.44
2:B:270:ILE:HG12	2:B:346:PHE:HB3	2.00	0.44
1:A:175:ASN:HB3	1:A:178:ILE:HD12	1.99	0.44
1:A:340:GLN:CB	1:A:351:THR:HG22	2.46	0.44
1:A:515:SER:CB	1:A:518:VAL:HG23	2.48	0.44
1:A:88:TRP:CH2	2:B:57:ASN:CB	3.01	0.44
1:A:98:ALA:HB2	1:A:349:LEU:O	2.18	0.44
1:A:163:SER:O	1:A:167:ILE:HG13	2.17	0.44
1:A:448:ARG:NE	1:A:473:THR:HB	2.30	0.44
2:B:378:GLU:O	2:B:382:ILE:HG13	2.17	0.44
1:A:278:GLN:NE2	1:A:281:LYS:HD2	2.33	0.44
1:A:143:ARG:HG2	1:A:143:ARG:NH1	2.30	0.44
1:A:502:ALA:O	1:A:506:ILE:HG13	2.18	0.44
1:A:500:GLN:HE21	1:A:500:GLN:HA	1.83	0.43
2:B:420:PRO:HB2	2:B:423:VAL:HG23	2.00	0.43
1:A:538:ALA:O	1:A:539:HIS:HB2	2.17	0.43
1:A:57:ASN:HA	1:A:129:ALA:O	2.18	0.43
1:A:451:LYS:O	1:A:471:ASP:N	2.52	0.43
1:A:248:GLU:OE1	1:A:248:GLU:HA	2.18	0.43
1:A:393:ILE:HB	1:A:423:VAL:CG1	2.43	0.43
2:B:97:PRO:C	2:B:99:GLY:N	2.77	0.43
1:A:228:LEU:HD22	1:A:242:GLN:OE1	2.19	0.42
2:B:169:GLU:OE2	2:B:169:GLU:HA	2.17	0.42
2:B:328:GLU:O	2:B:339:TYR:HA	2.19	0.42
1:A:112:GLY:O	1:A:116:PHE:N	2.51	0.42
1:A:318:TYR:CD1	4:A:999:GWE:H15A	2.54	0.42
2:B:60:VAL:HG22	2:B:61:PHE:N	2.34	0.42
1:A:335:GLY:HA2	1:A:367:GLN:OE1	2.19	0.42
1:A:317:VAL:HG23	1:A:349:LEU:HD23	2.02	0.42
2:B:23:GLN:HG2	2:B:133:PRO:HD3	2.02	0.42
2:B:46:LYS:HD3	2:B:116:PHE:CD2	2.54	0.42
2:B:365:VAL:O	2:B:369:THR:HG23	2.19	0.42
1:A:88:TRP:HA	1:A:88:TRP:CE3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:28:GLU:HA	2:B:135:ILE:HD11	2.00	0.42
1:A:116:PHE:CE1	1:A:164:MET:HE2	2.55	0.42
1:A:357:MET:HE2	1:A:374:LYS:NZ	2.35	0.42
1:A:438:GLU:OE1	1:A:459:THR:OG1	2.36	0.42
1:A:455:ALA:O	1:A:466:VAL:HA	2.20	0.42
1:A:78:ARG:O	1:A:82:LYS:HG3	2.20	0.42
1:A:226:PRO:HB3	1:A:235:HIS:CE1	2.54	0.42
2:B:332:GLN:NE2	2:B:424:LYS:HE2	2.35	0.41
1:A:59:PRO:HG2	1:A:76:ASP:HB3	2.02	0.41
2:B:253:THR:H	2:B:256:ASP:HB2	1.84	0.41
1:A:19:PRO:O	1:A:56:TYR:HA	2.20	0.41
1:A:94:ILE:HA	1:A:95:PRO:HD3	1.88	0.41
1:A:74:LEU:C	1:A:74:LEU:CD2	2.93	0.41
1:A:122:GLU:CD	1:A:122:GLU:N	2.75	0.41
2:B:353:LYS:CE	2:B:430:GLU:HB3	2.33	0.41
1:A:89:GLU:CD	1:A:89:GLU:H	2.29	0.41
1:A:483:TYR:CE2	1:A:487:GLN:NE2	2.89	0.41
2:B:97:PRO:C	2:B:99:GLY:H	2.28	0.41
1:A:101:LYS:O	4:A:999:GWE:H6	2.20	0.41
1:A:366:LYS:HZ1	2:B:394:GLN:HE21	1.68	0.41
1:A:498:ASP:HA	1:A:536:VAL:O	2.21	0.41
2:B:169:GLU:CB	2:B:170:PRO:HD3	2.51	0.41
2:B:342:TYR:CD1	2:B:342:TYR:C	2.98	0.41
1:A:224:GLU:HA	1:A:225:PRO:HD3	1.93	0.40
2:B:40:GLU:O	2:B:44:GLU:HG3	2.21	0.40
2:B:142:ILE:HD12	2:B:142:ILE:N	2.36	0.40
2:B:267:ALA:HB2	2:B:426:TRP:CZ3	2.56	0.40
1:A:52:PRO:O	1:A:53:GLU:C	2.65	0.40
1:A:116:PHE:CG	1:A:214:LEU:HD21	2.56	0.40
1:A:226:PRO:HB3	1:A:235:HIS:ND1	2.36	0.40
2:B:266:TRP:HZ3	2:B:426:TRP:CD1	2.39	0.40
2:B:342:TYR:HB3	2:B:348:ASN:HA	2.03	0.40
1:A:48:SER:O	1:A:50:ILE:HG23	2.22	0.40
1:A:93:GLY:O	2:B:137:ASN:HB3	2.22	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	525/560 (94%)	490 (93%)	26 (5%)	9 (2%)	7	5
2	B	392/440 (89%)	370 (94%)	20 (5%)	2 (0%)	24	27
All	All	917/1000 (92%)	860 (94%)	46 (5%)	11 (1%)	10	8

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	VAL
1	A	217	PRO
1	A	91	GLN
2	B	239	TRP
1	A	14	PRO
1	A	89	GLU
1	A	361	HIS
1	A	133	PRO
1	A	85	GLN
2	B	210	LEU
1	A	345	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	476/499 (95%)	449 (94%)	27 (6%)	18	23
2	B	363/400 (91%)	353 (97%)	10 (3%)	38	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	839/899 (93%)	802 (96%)	37 (4%)	25	34

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	14	PRO
1	A	53	GLU
1	A	89	GLU
1	A	94	ILE
1	A	122	GLU
1	A	138	GLU
1	A	143	ARG
1	A	160	PHE
1	A	205	LEU
1	A	210	LEU
1	A	214	LEU
1	A	217	PRO
1	A	230	MET
1	A	238	LYS
1	A	241	VAL
1	A	248	GLU
1	A	275	LYS
1	A	340	GLN
1	A	346	PHE
1	A	368	LEU
1	A	386	THR
1	A	402	TRP
1	A	404	GLU
1	A	423	VAL
1	A	459	THR
1	A	517	LEU
2	B	24	TRP
2	B	55	PRO
2	B	73	LYS
2	B	101	LYS
2	B	167	ILE
2	B	206	ARG
2	B	283	LEU
2	B	289	LEU
2	B	300	GLU
2	B	325	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	151	GLN
1	A	197	GLN
1	A	207	GLN
1	A	208	HIS
1	A	221	HIS
1	A	222	GLN
1	A	242	GLN
1	A	255	ASN
1	A	278	GLN
1	A	334	GLN
1	A	336	GLN
1	A	348	ASN
1	A	363	ASN
1	A	474	ASN
1	A	487	GLN
1	A	500	GLN
1	A	512	GLN
1	A	519	ASN
1	A	520	GLN
1	A	524	GLN
2	B	57	ASN
2	B	151	GLN
2	B	161	GLN
2	B	175	ASN
2	B	182	GLN
2	B	197	GLN
2	B	242	GLN
2	B	255	ASN
2	B	269	GLN
2	B	278	GLN
2	B	332	GLN
2	B	394	GLN
2	B	418	ASN
2	B	428	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CSD	A	280	1	4,7,8	0.74	0	1,8,10	0.38	0

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
1	CSD	A	280	1	-	1/2/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	280	CSD	N-CA-CB-SG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	GWE	A	999	-	34,38,38	2.86	12 (35%)	49,57,57	1.22	7 (14%)
3	PO4	A	1301	-	4,4,4	1.72	0	6,6,6	0.46	0
3	PO4	B	1300	-	4,4,4	1.76	0	6,6,6	0.45	0

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	GWE	A	999	-	-	2/25/29/29	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	GWE	C6-C5	6.91	1.50	1.38
4	A	999	GWE	C3-C4	6.20	1.48	1.38
4	A	999	GWE	C2-C1	6.11	1.51	1.40
4	A	999	GWE	C22-C21	6.04	1.48	1.38
4	A	999	GWE	C17-C18	5.38	1.51	1.40
4	A	999	GWE	C19-C20	4.75	1.50	1.39
4	A	999	GWE	C9-C10	3.30	1.43	1.37
4	A	999	GWE	C19-C18	-3.03	1.35	1.39
4	A	999	GWE	C11-C10	2.62	1.42	1.37
4	A	999	GWE	C8-C7	2.37	1.53	1.49
4	A	999	GWE	C9-C8	2.24	1.42	1.39
4	A	999	GWE	C13-C8	2.07	1.42	1.39

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	999	GWE	C2-C7-C8	2.76	123.97	119.56
4	A	999	GWE	C11-C10-C9	-2.67	120.25	123.50
4	A	999	GWE	C1-C2-C7	-2.18	119.48	122.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	999	GWE	C15-O2-C1	2.11	122.00	117.60
4	A	999	GWE	O2-C15-C16	-2.09	105.05	110.80
4	A	999	GWE	C22-C17-C18	-2.05	118.35	120.70
4	A	999	GWE	F2-C14-C12	-2.01	108.59	112.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

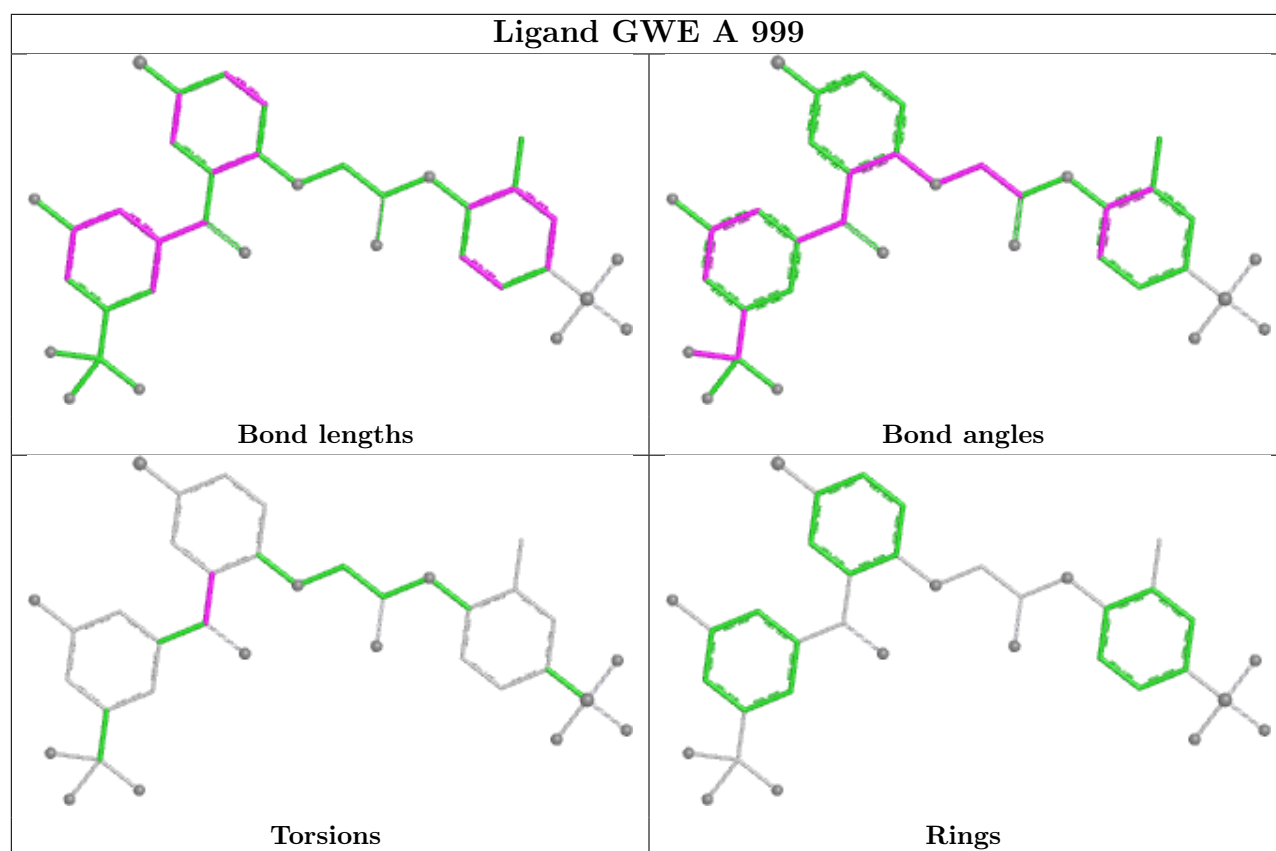
Mol	Chain	Res	Type	Atoms
4	A	999	GWE	C1-C2-C7-C8
4	A	999	GWE	C1-C2-C7-O1

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	999	GWE	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	531/560 (94%)	0.19	18 (3%) 48 45	28, 53, 95, 132	0
2	B	400/440 (90%)	0.18	14 (3%) 47 44	29, 53, 101, 127	0
All	All	931/1000 (93%)	0.19	32 (3%) 48 45	28, 53, 98, 132	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90	VAL	4.3
2	B	241	VAL	3.9
2	B	435	VAL	3.9
1	A	115	TYR	3.7
2	B	434	ILE	3.6
2	B	195	ILE	3.3
2	B	69	THR	3.3
1	A	24	TRP	3.3
1	A	116	PHE	3.2
1	A	91	GLN	3.1
2	B	95	PRO	3.1
1	A	359	GLY	3.0
1	A	52	PRO	3.0
2	B	437	ALA	2.9
2	B	240	THR	2.8
1	A	114	ALA	2.7
2	B	87	PHE	2.7
1	A	501	TYR	2.6
1	A	216	THR	2.5
1	A	21	VAL	2.4
1	A	92	LEU	2.4
2	B	210	LEU	2.4
1	A	135	ILE	2.4
2	B	116	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	212	TRP	2.4
1	A	112	GLY	2.3
1	A	217	PRO	2.2
2	B	177	ASP	2.2
1	A	245	VAL	2.1
1	A	539	HIS	2.1
2	B	104	LYS	2.1
1	A	113	ASP	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. 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
1	CSD	A	280	8/9	0.92	0.09	51,55,69,77	0

6.3 Carbohydrates [i](#)

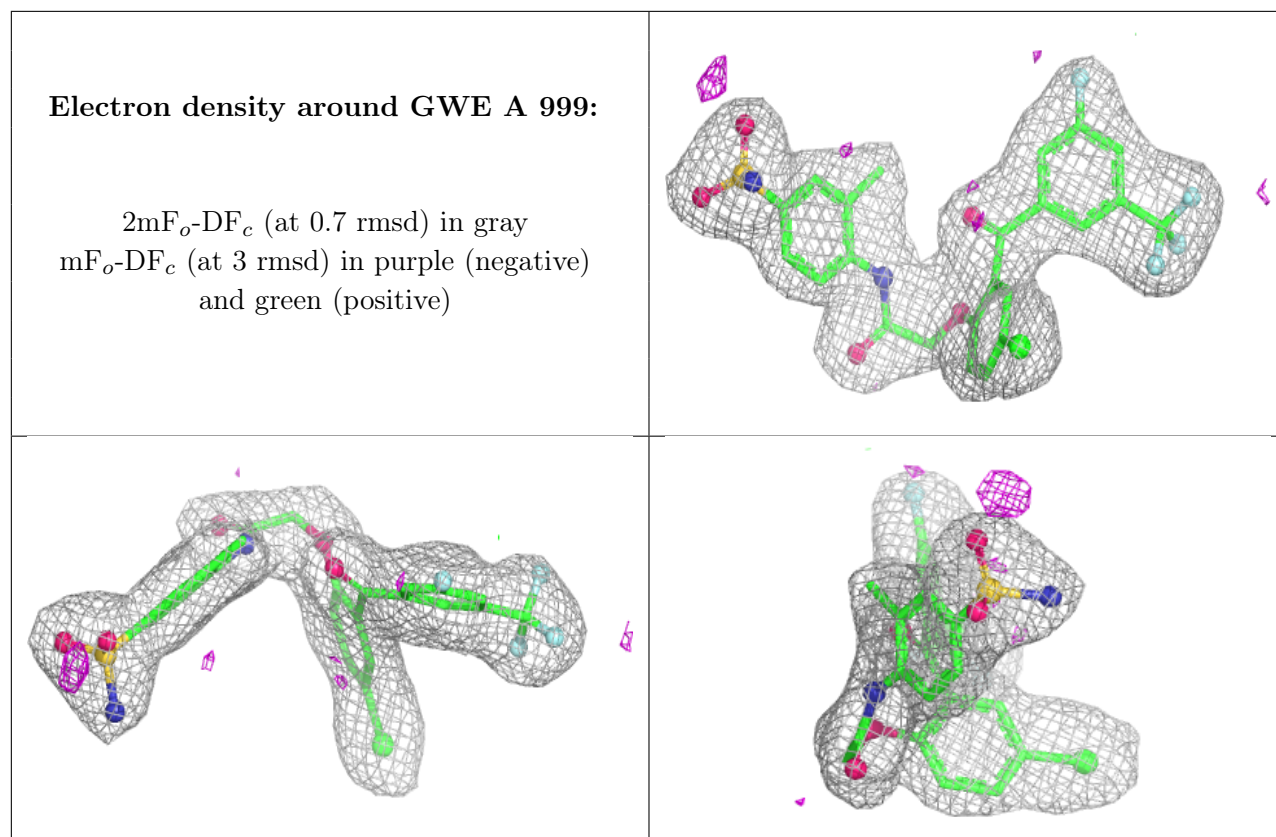
There are no oligosaccharides in this entry.

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	PO4	B	1300	5/5	0.72	0.11	117,123,129,132	0
3	PO4	A	1301	5/5	0.84	0.14	97,112,128,130	0
4	GWE	A	999	36/36	0.96	0.07	25,38,60,66	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.