



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:17 PM UTC

PDB ID : 5CEX / pdb_00005cex
Title : Crystal Structure of Fab 32H+109L, a putative precursor of the PGT121 family of potent HIV-1 antibodies
Authors : Garces, F.; Wilson, I.A.
Deposited on : 2015-07-07
Resolution : 2.10 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

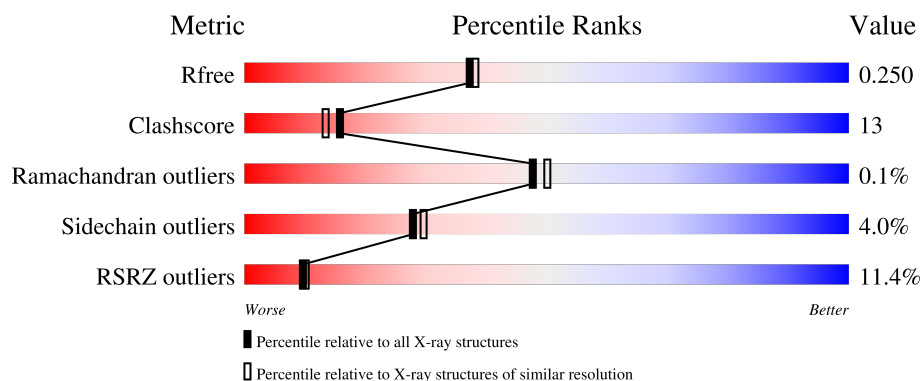
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	218	13% 73% 20% • 5%
1	C	218	10% 70% 23% • •
2	B	236	13% 79% 18% • •
2	D	236	8% 75% 20% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PG4	A	301	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7033 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.

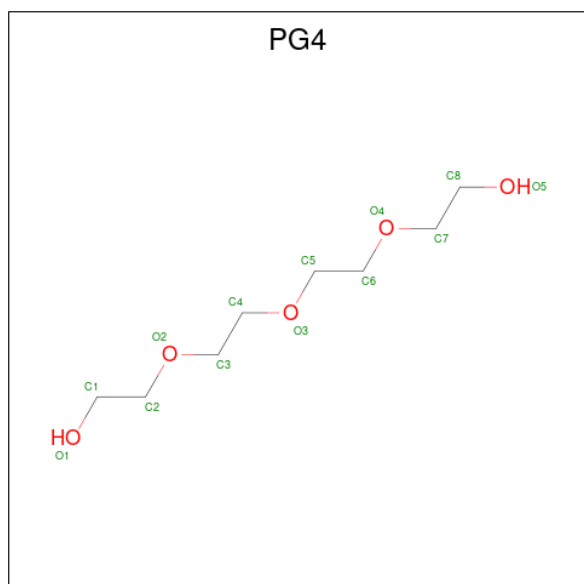
- Molecule 1 is a protein called Antibody 9H+3L Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	1	0
			1567	981	268	313	5			
1	C	213	Total	C	N	O	S	0	2	0
			1620	1015	276	324	5			

- Molecule 2 is a protein called Antibody 9H+3L Fab heavy chain.

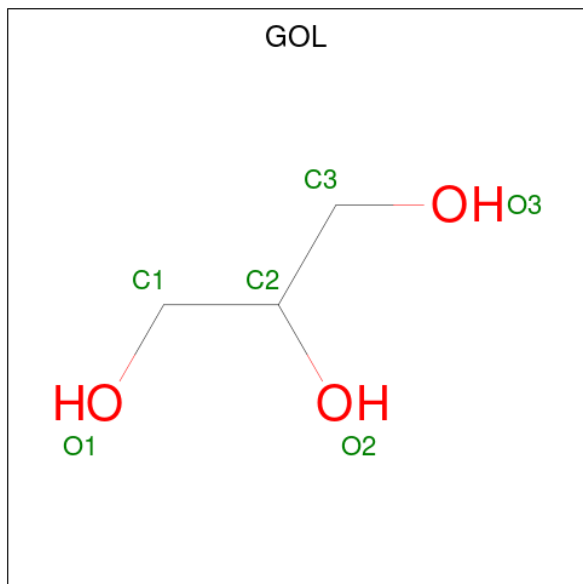
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	233	Total	C	N	O	S	0	0	0
			1761	1116	293	347	5			
2	D	232	Total	C	N	O	S	0	1	0
			1758	1116	292	345	5			

- Molecule 3 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



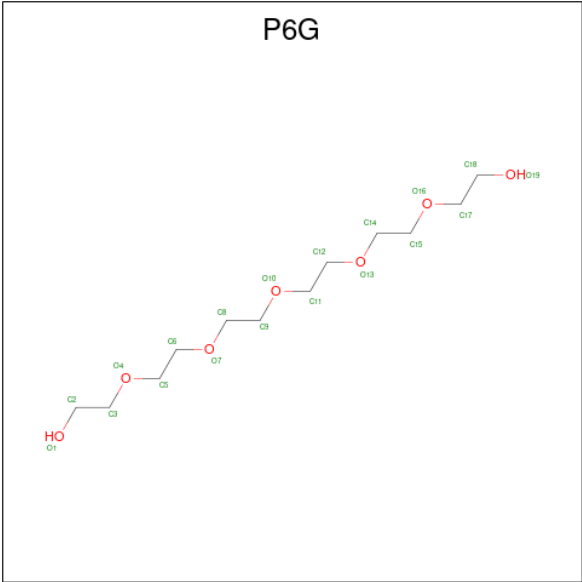
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	8	5		
3	C	1	Total	C	O	0	0
			13	8	5		
3	C	1	Total	C	O	0	0
			13	8	5		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



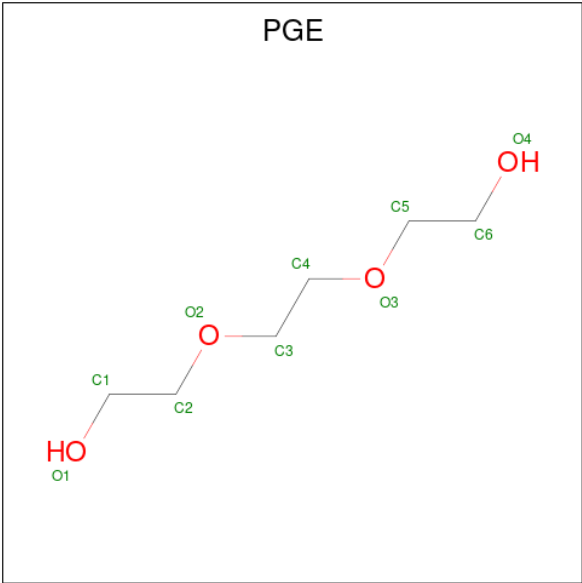
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: $C_{12}H_{26}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			19	12	7		
5	C	1	Total	C	O	0	0
			19	12	7		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is water.

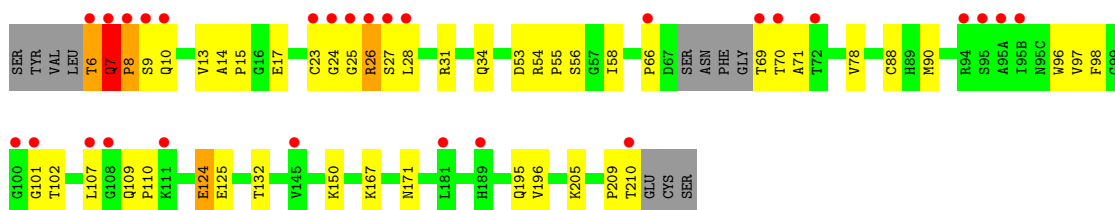
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	50	Total 50	O 50	0	0
7	B	46	Total 46	O 46	0	0
7	C	67	Total 67	O 67	0	0
7	D	53	Total 53	O 53	0	0

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.

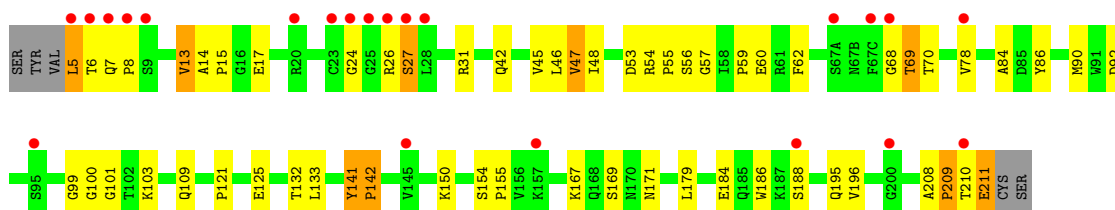
- Molecule 1: Antibody 9H+3L Fab light chain

Chain A: 



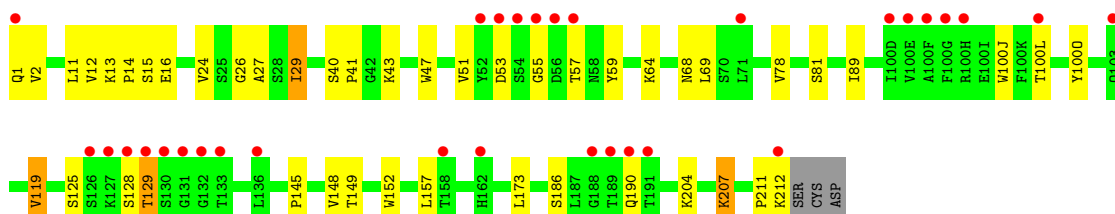
- Molecule 1: Antibody 9H+3L Fab light chain

Chain C: 



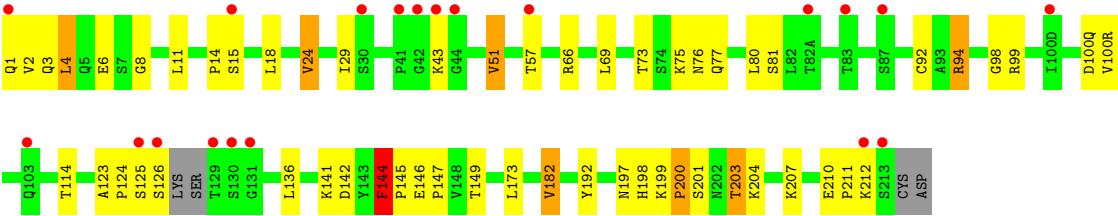
- Molecule 2: Antibody 9H+3L Fab heavy chain

Chain B: 



- Molecule 2: Antibody 9H+3L Fab heavy chain

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.95Å 108.73Å 72.70Å 90.00° 91.35° 90.00°	Depositor
Resolution (Å)	46.45 – 2.10 46.45 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.0 (46.45-2.10) 98.0 (46.45-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.200 , 0.244 0.220 , 0.250	Depositor DCC
R_{free} test set	3171 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.8	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7033	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, P6G, GOL, PGE

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	0.72	8/1609 (0.5%)	0.98	9/2197 (0.4%)
1	C	0.71	7/1664 (0.4%)	1.01	9/2274 (0.4%)
2	B	0.63	0/1808	0.87	0/2472
2	D	0.78	8/1807 (0.4%)	1.04	17/2468 (0.7%)
All	All	0.71	23/6888 (0.3%)	0.97	35/9411 (0.4%)

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 (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	124[A]	GLU	N-CA	7.33	1.55	1.46
1	A	124[B]	GLU	N-CA	7.33	1.55	1.46
1	C	142	PRO	N-CD	5.86	1.56	1.47
1	C	7	GLN	C-N	5.70	1.41	1.33
1	C	14	ALA	C-N	5.66	1.41	1.33
2	D	182	VAL	C-N	5.60	1.41	1.33
1	A	124[A]	GLU	CA-C	5.58	1.60	1.52
1	A	124[B]	GLU	CA-C	5.58	1.60	1.52
2	D	123	ALA	C-N	5.54	1.40	1.33
1	A	7	GLN	C-N	5.53	1.40	1.33
2	D	210	GLU	C-N	5.49	1.40	1.33
1	A	109	GLN	C-N	5.46	1.40	1.33
2	D	124	PRO	N-CD	5.26	1.55	1.47
1	C	208	ALA	C-N	5.24	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	211	PRO	N-CD	5.22	1.55	1.47
1	C	209	PRO	N-CD	5.21	1.55	1.47
2	D	147	PRO	N-CD	5.21	1.55	1.47
1	C	15	PRO	N-CD	5.20	1.55	1.47
2	D	200	PRO	N-CD	5.19	1.55	1.47
1	A	110	PRO	N-CD	5.16	1.54	1.47
1	A	8	PRO	N-CD	5.13	1.54	1.47
2	D	4	LEU	C-O	-5.12	1.17	1.24
1	C	8	PRO	N-CD	5.06	1.54	1.47

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	71	ALA	O-C-N	-11.32	108.94	122.87
2	D	146	GLU	C-N-CD	8.55	139.40	120.60
1	C	141	TYR	C-N-CD	7.76	137.66	120.60
1	C	7	GLN	CA-C-N	-6.87	112.69	119.76
1	C	7	GLN	C-N-CA	-6.87	112.69	119.76
1	A	7	GLN	CA-C-N	-6.70	113.08	119.85
1	A	7	GLN	C-N-CA	-6.70	113.08	119.85
1	C	69	THR	N-CA-C	6.60	119.33	110.35
1	A	109	GLN	CA-C-N	-6.31	113.45	119.76
1	A	109	GLN	C-N-CA	-6.31	113.45	119.76
2	D	144	PHE	C-N-CD	6.29	134.44	120.60
1	C	195	GLN	CA-C-N	-6.29	115.39	123.19
1	C	195	GLN	C-N-CA	-6.29	115.39	123.19
1	A	96	TRP	CA-C-N	-6.27	115.41	123.19
1	A	96	TRP	C-N-CA	-6.27	115.41	123.19
2	D	182	VAL	CA-C-N	-6.27	113.49	119.76
2	D	182	VAL	C-N-CA	-6.27	113.49	119.76
2	D	210	GLU	CA-C-N	-6.22	113.36	119.76
2	D	210	GLU	C-N-CA	-6.22	113.36	119.76
2	D	144	PHE	CA-C-N	-6.20	112.11	127.00
2	D	144	PHE	C-N-CA	-6.20	112.11	127.00
1	C	14	ALA	CA-C-N	-6.18	113.40	119.76
1	C	14	ALA	C-N-CA	-6.18	113.40	119.76
2	D	8	GLY	CA-C-N	6.15	125.77	119.56
2	D	8	GLY	C-N-CA	6.15	125.77	119.56
2	D	123	ALA	CA-C-N	-6.10	113.69	119.85
2	D	123	ALA	C-N-CA	-6.10	113.69	119.85
2	D	144	PHE	CB-CA-C	5.80	119.24	109.32
1	C	100	GLY	N-CA-C	5.51	120.46	113.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	GLN	N-CA-C	5.49	117.74	109.23
2	D	147	PRO	CA-N-CD	-5.38	103.97	111.50
2	D	136	LEU	CA-CB-CG	5.35	135.03	116.30
1	A	97	VAL	N-CA-C	5.26	115.53	108.17
2	D	199	LYS	CA-C-N	-5.02	113.48	119.05
2	D	199	LYS	C-N-CA	-5.02	113.48	119.05

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	144	PHE	Peptide

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	1567	0	1516	44	0
1	C	1620	0	1571	68	0
2	B	1761	0	1717	46	0
2	D	1758	0	1719	36	0
3	A	13	0	18	11	0
3	C	26	0	36	11	0
4	B	6	0	8	0	0
4	C	18	0	24	5	0
5	B	19	0	26	4	0
5	C	19	0	26	7	0
6	C	10	0	14	1	0
7	A	50	0	0	1	0
7	B	46	0	0	1	0
7	C	67	0	0	0	0
7	D	53	0	0	2	0
All	All	7033	0	6675	179	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 13.

All (179) 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:1:GLN:HE21	1:C:57:GLY:CA	1.55	1.17
1:A:7:GLN:HG3	1:A:101:GLY:H	1.11	1.14
2:B:1:GLN:NE2	1:C:57:GLY:HA2	1.66	1.08
2:B:1:GLN:NE2	1:C:57:GLY:CA	2.21	1.04
2:B:1:GLN:HE21	1:C:57:GLY:C	1.65	1.03
2:D:94:ARG:NH1	7:D:301:HOH:O	1.93	1.00
2:B:125:SER:O	2:B:129:THR:HG23	1.62	1.00
1:A:15:PRO:HD3	1:A:107:LEU:O	1.64	0.97
1:C:42:GLN:HE21	5:C:304:P6G:H31	1.28	0.96
1:C:45:VAL:HG13	3:C:305:PG4:H11	1.45	0.95
1:C:6:THR:O	1:C:24:GLY:HA2	1.69	0.92
1:A:7:GLN:HG3	1:A:101:GLY:N	1.85	0.90
1:A:7:GLN:CG	1:A:101:GLY:H	1.86	0.89
2:B:1:GLN:NE2	1:C:57:GLY:O	2.06	0.88
2:B:1:GLN:HG2	1:C:57:GLY:CA	2.05	0.85
1:C:27:SER:OG	1:C:92:ASP:OD1	1.95	0.85
1:A:54:ARG:HH12	3:A:301:PG4:H12	1.40	0.85
1:A:54:ARG:H	3:A:301:PG4:H52	1.43	0.82
2:B:1:GLN:HG2	1:C:57:GLY:C	2.04	0.82
1:C:13:VAL:CG1	1:C:78:VAL:HG21	2.10	0.82
2:B:1:GLN:CD	1:C:57:GLY:HA2	2.07	0.79
2:B:1:GLN:NE2	1:C:57:GLY:C	2.41	0.77
1:C:53:ASP:HA	4:C:303:GOL:H31	1.66	0.76
2:D:201:SER:OG	2:D:203:THR:HG23	1.85	0.76
1:A:7:GLN:NE2	1:A:88:CYS:SG	2.60	0.74
2:B:2:VAL:HG12	2:B:27:ALA:HB2	1.67	0.74
2:B:1:GLN:CG	1:C:57:GLY:C	2.59	0.74
3:A:301:PG4:H62	2:D:99:ARG:H	1.52	0.74
1:C:209:PRO:O	1:C:211:GLU:HG3	1.88	0.73
2:B:125:SER:N	2:B:128:SER:OG	2.12	0.72
1:C:133[B]:LEU:HD12	1:C:179:LEU:HD23	1.71	0.71
2:D:142:ASP:HB3	2:D:173:LEU:HD13	1.73	0.71
1:A:31:ARG:NH2	1:A:66:PRO:O	2.23	0.71
1:A:7:GLN:OE1	1:A:101:GLY:HA2	1.92	0.70
2:B:89:ILE:CD1	5:B:302:P6G:H21	2.21	0.69
1:A:125:GLU:OE1	7:A:401:HOH:O	2.10	0.69
2:D:2:VAL:HG23	2:D:100(R):VAL:HG11	1.74	0.69
1:C:5:LEU:HD11	1:C:90:MET:HB3	1.74	0.69
3:A:301:PG4:H62	2:D:98:GLY:HA2	1.74	0.69
2:D:51:VAL:HG12	2:D:57:THR:HG22	1.73	0.68
1:C:57:GLY:H	3:C:306:PG4:H12	1.59	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:GLN:HG2	1:C:57:GLY:HA2	1.75	0.68
1:C:46:LEU:O	3:C:305:PG4:H22	1.93	0.67
1:C:6:THR:O	1:C:24:GLY:CA	2.43	0.67
2:B:186:SER:HB2	2:B:190:GLN:HG3	1.77	0.66
1:C:59:PRO:HB3	4:C:302:GOL:H32	1.76	0.66
1:C:42:GLN:HE22	5:C:304:P6G:H92	1.60	0.65
2:B:89:ILE:HD13	5:B:302:P6G:H21	1.78	0.65
1:C:184:GLU:O	1:C:188:SER:OG	2.13	0.65
2:B:1:GLN:CG	1:C:57:GLY:HA2	2.27	0.64
1:A:54:ARG:N	3:A:301:PG4:H52	2.13	0.63
1:A:27:SER:O	1:A:69:THR:HG22	1.97	0.63
3:C:305:PG4:H52	2:D:100(Q):ASP:OD2	1.99	0.63
2:D:2:VAL:CG2	2:D:100(R):VAL:HG21	2.29	0.63
2:B:51:VAL:HG23	2:B:57:THR:HG22	1.81	0.62
1:C:5:LEU:HB3	1:C:99:GLY:HA2	1.81	0.62
6:C:307:PGE:O1	6:C:307:PGE:H3	1.98	0.62
2:B:100(J):TRP:HZ3	2:B:100(L):THR:HG23	1.64	0.62
2:B:1:GLN:CG	1:C:57:GLY:CA	2.78	0.62
1:C:13:VAL:HG11	1:C:78:VAL:HG21	1.81	0.61
1:C:42:GLN:HE21	5:C:304:P6G:C3	2.09	0.61
1:C:13:VAL:HG13	1:C:17:GLU:HG2	1.82	0.61
1:C:210:THR:C	1:C:211:GLU:HG3	2.24	0.61
1:A:209:PRO:O	1:A:210:THR:HG23	2.01	0.61
2:B:119:VAL:HG23	2:B:207:LYS:HG3	1.82	0.61
1:A:53:ASP:OD1	3:A:301:PG4:H61	2.01	0.60
1:C:13:VAL:HG12	1:C:78:VAL:HG21	1.84	0.60
2:B:125:SER:H	2:B:128:SER:HG	1.48	0.60
1:A:69:THR:O	1:A:70:THR:HG22	2.01	0.60
1:A:70:THR:O	1:A:70:THR:HG23	2.01	0.60
3:A:301:PG4:H62	2:D:99:ARG:N	2.15	0.59
2:B:11:LEU:HB2	2:B:145:PRO:HG3	1.83	0.59
1:C:31:ARG:HH12	1:C:69:THR:HB	1.68	0.58
2:D:197:ASN:OD1	2:D:204:LYS:HE2	2.03	0.58
1:C:57:GLY:N	3:C:306:PG4:H12	2.18	0.58
2:D:2:VAL:HG21	2:D:100(R):VAL:HG21	1.86	0.58
1:C:5:LEU:HD11	1:C:90:MET:CB	2.34	0.58
1:C:42:GLN:NE2	5:C:304:P6G:H31	2.09	0.57
1:A:53:ASP:OD1	2:D:99:ARG:HD3	2.04	0.57
2:D:3:GLN:C	2:D:4:LEU:HD23	2.29	0.57
2:D:149[B]:THR:OG1	2:D:197:ASN:HB3	2.05	0.57
1:A:27:SER:HB2	1:A:90:MET:SD	2.45	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:125:SER:O	2:B:128:SER:N	2.38	0.56
2:D:66:ARG:NH2	7:D:303:HOH:O	2.38	0.56
1:A:125:GLU:OE2	1:A:132:THR:HG23	2.06	0.55
1:C:26:ARG:HD2	1:C:68:GLY:O	2.04	0.55
2:D:182:VAL:HG11	2:D:192:TYR:OH	2.06	0.55
1:C:141:TYR:CG	1:C:142:PRO:HA	2.43	0.54
1:A:167:LYS:HE2	1:A:171:ASN:HA	1.90	0.54
3:A:301:PG4:H62	2:D:98:GLY:CA	2.37	0.54
1:C:121:PRO:HD3	1:C:133[B]:LEU:HD23	1.90	0.54
2:D:6:GLU:HG3	2:D:92:CYS:SG	2.48	0.54
1:A:15:PRO:CD	1:A:107:LEU:O	2.50	0.53
2:D:198:HIS:CD2	2:D:200:PRO:HD2	2.43	0.53
2:D:75:LYS:HB2	2:D:77:GLN:HG2	1.89	0.53
2:B:2:VAL:HG13	7:B:425:HOH:O	2.09	0.52
2:D:24:VAL:HG22	2:D:76:ASN:ND2	2.26	0.51
2:D:125:SER:O	2:D:126:SER:HB3	2.10	0.51
2:B:152:TRP:HB3	2:B:157:LEU:HD23	1.91	0.51
2:B:211:PRO:O	2:B:212:LYS:HB2	2.10	0.51
1:A:56:SER:O	3:C:306:PG4:H72	2.11	0.51
1:A:13:VAL:CG1	1:A:78:VAL:HG21	2.40	0.50
2:B:40:SER:HB2	2:B:43:LYS:HG3	1.93	0.50
2:B:2:VAL:HG12	2:B:27:ALA:CB	2.39	0.50
3:A:301:PG4:C6	2:D:99:ARG:H	2.23	0.49
1:A:53:ASP:HA	3:A:301:PG4:H42	1.94	0.49
1:C:210:THR:C	1:C:211:GLU:CG	2.85	0.49
1:C:56:SER:OG	3:C:306:PG4:H32	2.12	0.49
1:A:98:PHE:HE1	2:B:47:TRP:HB2	1.78	0.48
1:C:47:VAL:HG12	1:C:48:ILE:HG12	1.93	0.48
1:A:209:PRO:O	1:A:210:THR:CG2	2.61	0.48
1:C:55:PRO:HB2	3:C:305:PG4:H42	1.96	0.47
2:D:204:LYS:HE3	2:D:204:LYS:HB2	1.61	0.47
1:C:69:THR:HG22	1:C:70:THR:N	2.28	0.47
2:D:144:PHE:HA	2:D:145:PRO:HA	1.67	0.47
1:A:23:CYS:SG	1:A:24:GLY:N	2.88	0.46
1:C:103:LYS:HG3	4:C:301:GOL:H32	1.97	0.46
1:A:27:SER:HB3	1:A:69:THR:N	2.31	0.46
1:A:150:LYS:HE2	1:A:195:GLN:OE1	2.15	0.46
1:C:109:GLN:HB2	1:C:141:TYR:CE1	2.50	0.46
5:C:304:P6G:H92	5:C:304:P6G:H62	1.44	0.45
1:C:169:SER:HA	5:C:304:P6G:H61	1.97	0.45
1:C:59:PRO:CB	4:C:302:GOL:H32	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:GLU:OE2	1:C:132:THR:HG23	2.17	0.45
1:A:8:PRO:O	1:A:102:THR:HA	2.17	0.45
2:B:29:ILE:HA	2:B:29:ILE:HD12	1.76	0.45
1:C:186:TRP:O	1:C:209:PRO:HG2	2.17	0.44
1:A:6:THR:CB	1:A:24:GLY:HA3	2.47	0.44
1:A:6:THR:HB	1:A:24:GLY:HA3	1.99	0.44
1:A:13:VAL:HG12	1:A:78:VAL:HG21	1.98	0.44
2:B:26:GLY:O	1:C:60:GLU:HG3	2.16	0.44
1:C:42:GLN:NE2	5:C:304:P6G:H62	2.32	0.44
1:C:54:ARG:NE	1:C:62:PHE:O	2.47	0.44
1:C:26:ARG:O	1:C:27:SER:HB3	2.18	0.44
1:A:14:ALA:O	1:A:17:GLU:HG2	2.17	0.44
1:A:209:PRO:C	1:A:210:THR:HG23	2.43	0.44
2:B:53:ASP:C	2:B:55:GLY:H	2.26	0.44
2:B:68:ASN:HB3	2:B:81:SER:OG	2.18	0.43
2:D:11:LEU:HD22	2:D:114:THR:HG22	2.00	0.43
2:B:53:ASP:C	2:B:55:GLY:N	2.76	0.43
2:B:204:LYS:HE3	2:B:204:LYS:HB2	1.80	0.43
1:C:5:LEU:HB3	1:C:99:GLY:CA	2.47	0.43
1:C:121:PRO:HD3	1:C:133[B]:LEU:CD2	2.49	0.43
2:D:182:VAL:CG1	2:D:192:TYR:OH	2.66	0.43
3:A:301:PG4:H71	2:D:98:GLY:HA3	2.01	0.43
2:B:100(J):TRP:HZ3	2:B:100(L):THR:CG2	2.32	0.42
1:C:141:TYR:HA	1:C:142:PRO:C	2.44	0.42
1:C:56:SER:HA	3:C:306:PG4:H32	2.01	0.42
1:C:150:LYS:HD3	1:C:155:PRO:HA	2.01	0.42
1:A:17:GLU:O	1:A:78:VAL:HG22	2.19	0.42
1:C:27:SER:OG	1:C:92:ASP:CG	2.61	0.42
2:D:4:LEU:HD23	2:D:4:LEU:N	2.34	0.42
2:B:29:ILE:HD11	2:B:78:VAL:HG23	2.02	0.42
2:D:18:LEU:O	2:D:81:SER:HA	2.20	0.42
2:D:69:LEU:CD2	2:D:80:LEU:HD13	2.49	0.42
1:A:55:PRO:HD2	1:A:58:ILE:HG13	2.01	0.42
1:A:56:SER:O	3:C:305:PG4:H72	2.20	0.42
2:B:41:PRO:CG	5:B:302:P6G:H152	2.49	0.42
2:B:89:ILE:HD12	5:B:302:P6G:H21	1.99	0.42
2:D:14:PRO:O	2:D:15:SER:HB2	2.20	0.42
2:D:29:ILE:HD11	2:D:73:THR:HA	2.02	0.42
2:D:144:PHE:O	2:D:198:HIS:CE1	2.73	0.42
1:C:154:SER:HA	1:C:155:PRO:HD3	1.88	0.41
1:A:25:GLY:C	1:A:26:ARG:HG3	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:PRO:HB3	3:C:305:PG4:H62	2.03	0.41
2:D:182:VAL:HG11	2:D:192:TYR:CZ	2.55	0.41
1:A:7:GLN:OE1	1:A:101:GLY:CA	2.67	0.41
1:A:34:GLN:HG2	2:B:100(O):TYR:HB3	2.03	0.41
2:B:14:PRO:O	2:B:15:SER:HB2	2.21	0.41
1:C:84:ALA:HA	4:C:301:GOL:H12	2.02	0.41
1:A:28:LEU:HD12	1:A:28:LEU:HA	1.96	0.41
2:B:12:VAL:HG13	2:B:16:GLU:OE1	2.20	0.41
1:A:7:GLN:HG3	1:A:101:GLY:CA	2.50	0.41
1:C:167:LYS:HE2	1:C:171:ASN:HA	2.02	0.41
2:B:125:SER:C	2:B:129:THR:HG23	2.40	0.40
2:B:59:TYR:HE1	2:B:69:LEU:HG	1.86	0.40
1:C:17:GLU:O	1:C:78:VAL:HG22	2.20	0.40
1:C:86:TYR:O	1:C:101:GLY:HA2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	204/218 (94%)	198 (97%)	6 (3%)	0	100	100
1	C	213/218 (98%)	210 (99%)	2 (1%)	1 (0%)	24	22
2	B	231/236 (98%)	228 (99%)	3 (1%)	0	100	100
2	D	229/236 (97%)	224 (98%)	5 (2%)	0	100	100
All	All	877/908 (97%)	860 (98%)	16 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	27	SER

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	173/183 (94%)	167 (96%)	6 (4%)	32	35
1	C	180/183 (98%)	175 (97%)	5 (3%)	38	43
2	B	200/204 (98%)	190 (95%)	10 (5%)	22	22
2	D	201/204 (98%)	192 (96%)	9 (4%)	24	25
All	All	754/774 (97%)	724 (96%)	30 (4%)	28	29

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	7	GLN
1	A	9	SER
1	A	26	ARG
1	A	196	VAL
1	A	205	LYS
2	B	13	LYS
2	B	24	VAL
2	B	29	ILE
2	B	64	LYS
2	B	119	VAL
2	B	129	THR
2	B	148	VAL
2	B	149	THR
2	B	173	LEU
2	B	207	LYS
1	C	5	LEU
1	C	13	VAL
1	C	47	VAL
1	C	196	VAL
1	C	211	GLU
2	D	1	GLN
2	D	24	VAL
2	D	43	LYS

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Mol	Chain	Res	Type
2	D	51	VAL
2	D	94	ARG
2	D	141	LYS
2	D	203	THR
2	D	207	LYS
2	D	212	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	1	GLN
2	B	96	GLN
1	C	38	GLN
1	C	42	GLN
2	D	5	GLN
2	D	32	HIS
2	D	39	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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	$\# Z > 2$
3	PG4	C	306	-	12,12,12	0.40	0	11,11,11	0.22	0
4	GOL	C	303	-	5,5,5	0.24	0	5,5,5	0.35	0
3	PG4	C	305	-	12,12,12	0.86	0	11,11,11	0.44	0
3	PG4	A	301	-	12,12,12	0.84	0	11,11,11	0.73	0
4	GOL	B	301	-	5,5,5	0.39	0	5,5,5	0.69	0
4	GOL	C	302	-	5,5,5	0.18	0	5,5,5	0.55	0
5	P6G	B	302	-	18,18,18	0.46	0	17,17,17	0.25	0
5	P6G	C	304	-	18,18,18	0.80	0	17,17,17	0.41	0
6	PGE	C	307	-	9,9,9	0.28	0	8,8,8	0.59	0
4	GOL	C	301	-	5,5,5	0.21	0	5,5,5	0.28	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
3	PG4	C	306	-	-	6/10/10/10	-
4	GOL	C	303	-	-	3/4/4/4	-
3	PG4	C	305	-	-	6/10/10/10	-
3	PG4	A	301	-	-	5/10/10/10	-
4	GOL	B	301	-	-	2/4/4/4	-
4	GOL	C	302	-	-	0/4/4/4	-
5	P6G	B	302	-	-	9/16/16/16	-
5	P6G	C	304	-	-	7/16/16/16	-
6	PGE	C	307	-	-	4/7/7/7	-
4	GOL	C	301	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	GOL	C1-C2-C3-O3
4	C	301	GOL	C1-C2-C3-O3

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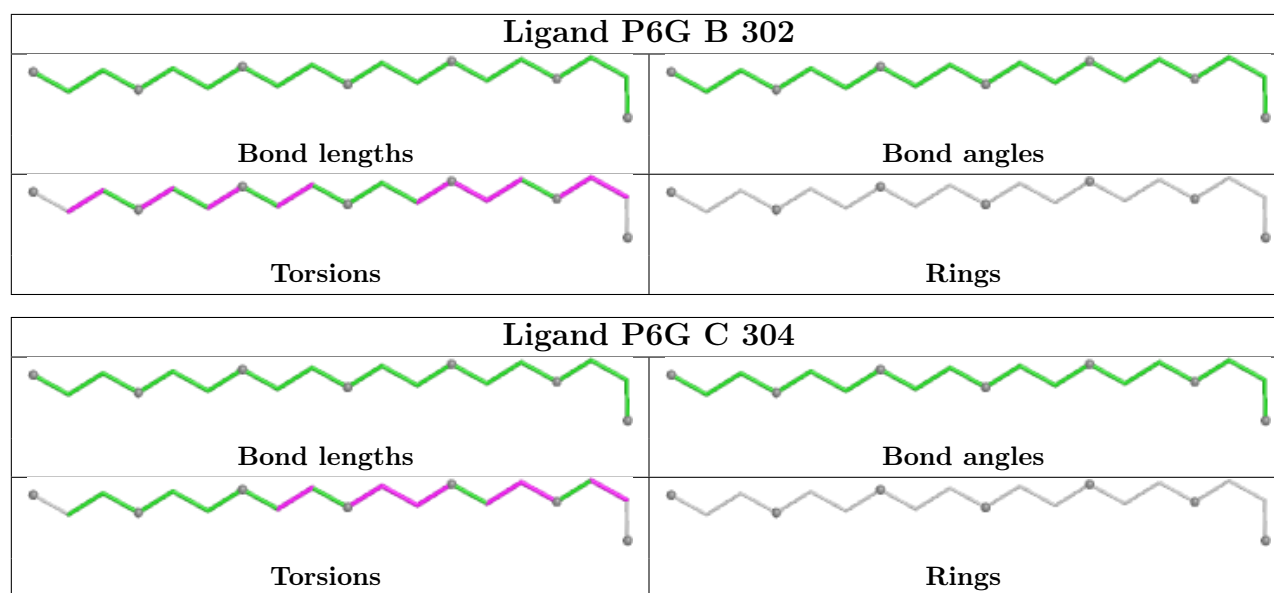
Mol	Chain	Res	Type	Atoms
5	C	304	P6G	C9-C8-O7-C6
3	C	305	PG4	C6-C5-O3-C4
3	C	306	PG4	O2-C3-C4-O3
3	C	306	PG4	O3-C5-C6-O4
5	C	304	P6G	O7-C8-C9-O10
3	C	305	PG4	C5-C6-O4-C7
3	C	305	PG4	O3-C5-C6-O4
3	C	305	PG4	O4-C7-C8-O5
5	B	302	P6G	O1-C2-C3-O4
5	C	304	P6G	O10-C11-C12-O13
3	A	301	PG4	O2-C3-C4-O3
4	C	303	GOL	O1-C1-C2-C3
4	B	301	GOL	O2-C2-C3-O3
4	C	301	GOL	O2-C2-C3-O3
5	C	304	P6G	O1-C2-C3-O4
5	B	302	P6G	O4-C5-C6-O7
6	C	307	PGE	C1-C2-O2-C3
6	C	307	PGE	O2-C3-C4-O3
5	C	304	P6G	O4-C5-C6-O7
4	C	303	GOL	O1-C1-C2-O2
3	A	301	PG4	O4-C7-C8-O5
6	C	307	PGE	O3-C5-C6-O4
5	B	302	P6G	C14-C15-O16-C17
3	C	306	PG4	C3-C4-O3-C5
3	C	306	PG4	O4-C7-C8-O5
3	C	305	PG4	C8-C7-O4-C6
5	B	302	P6G	C9-C8-O7-C6
3	C	306	PG4	C5-C6-O4-C7
5	B	302	P6G	C15-C14-O13-C12
3	C	306	PG4	O1-C1-C2-O2
3	A	301	PG4	C8-C7-O4-C6
5	B	302	P6G	O10-C11-C12-O13
5	B	302	P6G	C5-C6-O7-C8
3	A	301	PG4	C3-C4-O3-C5
4	C	303	GOL	O2-C2-C3-O3
3	A	301	PG4	C1-C2-O2-C3
5	C	304	P6G	C8-C9-O10-C11
6	C	307	PGE	C3-C4-O3-C5
5	B	302	P6G	O16-C17-C18-O19
5	C	304	P6G	C6-C5-O4-C3
5	B	302	P6G	C2-C3-O4-C5
3	C	305	PG4	O2-C3-C4-O3

There are no ring outliers.

9 monomers are involved in 39 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	306	PG4	5	0
4	C	303	GOL	1	0
3	C	305	PG4	6	0
3	A	301	PG4	11	0
4	C	302	GOL	2	0
5	B	302	P6G	4	0
5	C	304	P6G	7	0
6	C	307	PGE	1	0
4	C	301	GOL	2	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	207/218 (94%)	0.96	28 (13%) 7 7	19, 37, 70, 96	2 (0%)
1	C	213/218 (97%)	0.62	22 (10%) 12 12	15, 33, 57, 84	4 (1%)
2	B	233/236 (98%)	0.68	31 (13%) 7 7	24, 37, 66, 84	0
2	D	232/236 (98%)	0.78	20 (8%) 16 17	16, 35, 57, 83	5 (2%)
All	All	885/908 (97%)	0.76	101 (11%) 10 10	15, 36, 65, 96	11 (1%)

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	72	THR	10.9
2	D	87	SER	10.7
2	D	30	SER	9.8
1	A	24	GLY	7.2
1	C	25	GLY	7.1
1	A	6	THR	7.0
2	D	83	THR	6.8
2	D	213	SER	6.3
1	A	107	LEU	6.0
2	D	57	THR	6.0
1	A	69	THR	5.9
1	A	95	SER	5.7
1	C	210	THR	5.6
1	C	5	LEU	5.3
2	D	129	THR	5.2
1	C	8	PRO	4.9
1	C	20	ARG	4.9
2	B	1	GLN	4.7
1	A	95(A)	ALA	4.7
1	A	7	GLN	4.7
1	C	78	VAL	4.5

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Mol	Chain	Res	Type	RSRZ
2	D	42	GLY	4.5
1	A	8	PRO	4.4
2	B	100(F)	ALA	4.4
1	A	210	THR	4.3
2	B	127	LYS	4.3
1	A	28	LEU	4.1
1	A	9	SER	4.0
2	B	100(G)	PHE	3.9
1	C	68	GLY	3.6
2	B	100(E)	VAL	3.6
1	A	70	THR	3.6
2	B	131	GLY	3.5
1	A	27	SER	3.3
2	D	130	SER	3.3
2	B	53	ASP	3.3
1	C	27	SER	3.3
2	B	132	GLY	3.2
1	A	95(B)	ILE	3.1
2	D	41	PRO	3.0
2	D	103	GLN	3.0
2	D	100(D)	ILE	3.0
2	B	212	LYS	2.9
2	D	126	SER	2.9
2	D	131	GLY	2.9
2	B	54	SER	2.9
2	B	126	SER	2.8
1	C	26	ARG	2.8
1	C	24	GLY	2.8
1	A	94	ARG	2.8
1	C	67(A)	SER	2.8
2	B	103	GLN	2.8
1	C	67(C)	PHE	2.7
1	C	28	LEU	2.7
2	B	100(L)	THR	2.7
2	B	129	THR	2.7
1	C	7	GLN	2.7
2	D	125	SER	2.7
1	A	10	GLN	2.6
2	B	190	GLN	2.6
2	B	100(D)	ILE	2.6
2	D	44	GLY	2.6
1	C	95	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	128	SER	2.6
2	B	100(H)	ARG	2.6
1	A	101	GLY	2.6
1	C	23	CYS	2.6
1	A	145	VAL	2.5
1	A	111	LYS	2.5
1	A	25	GLY	2.5
1	A	26	ARG	2.5
2	D	15	SER	2.5
2	B	55	GLY	2.5
2	B	133	THR	2.4
1	C	157	LYS	2.4
2	B	188	GLY	2.4
2	B	158	THR	2.4
2	D	212	LYS	2.4
2	B	162	HIS	2.4
2	B	52	TYR	2.3
1	A	108	GLY	2.3
2	B	130	SER	2.3
1	C	145	VAL	2.3
1	C	6	THR	2.3
1	A	23	CYS	2.3
1	C	9	SER	2.2
2	B	191	THR	2.2
2	B	71	LEU	2.2
1	A	189	HIS	2.2
2	D	43	LYS	2.2
1	A	100	GLY	2.2
2	B	136	LEU	2.2
2	D	1	GLN	2.1
1	A	66	PRO	2.1
2	B	57	THR	2.1
1	C	200	GLY	2.1
2	B	189	THR	2.1
2	B	56	ASP	2.1
2	D	82(A)	THR	2.1
1	C	188	SER	2.1
1	A	181	LEU	2.0

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

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](#)

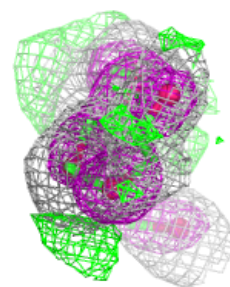
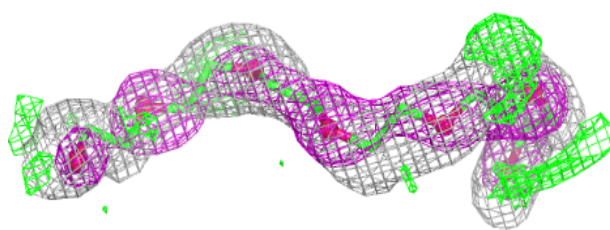
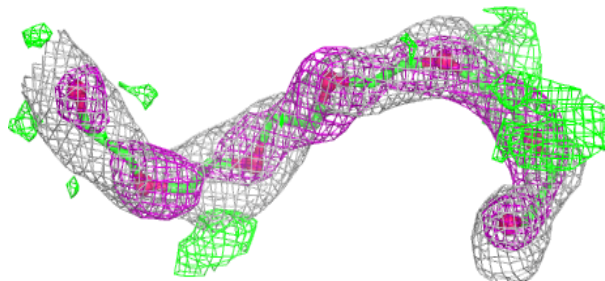
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	P6G	B	302	19/19	0.66	0.17	20,20,20,20	0
6	PGE	C	307	10/10	0.69	0.22	20,20,20,20	0
4	GOL	C	303	6/6	0.73	0.20	20,20,20,20	0
3	PG4	C	306	13/13	0.76	0.17	43,50,56,63	0
3	PG4	C	305	13/13	0.77	0.19	28,44,57,62	0
3	PG4	A	301	13/13	0.79	0.17	36,42,53,56	0
4	GOL	B	301	6/6	0.80	0.20	38,42,49,63	0
4	GOL	C	302	6/6	0.83	0.15	44,52,55,58	0
4	GOL	C	301	6/6	0.89	0.13	37,43,56,58	0
5	P6G	C	304	19/19	0.92	0.11	29,41,47,55	0

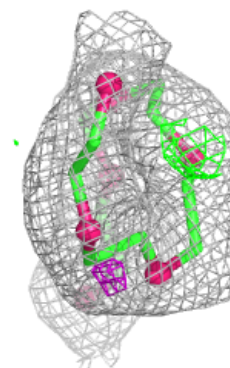
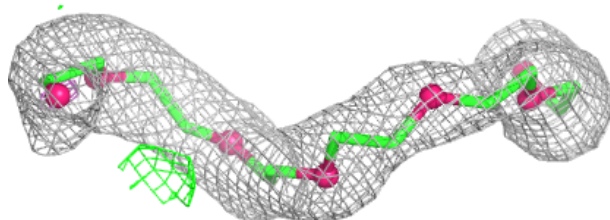
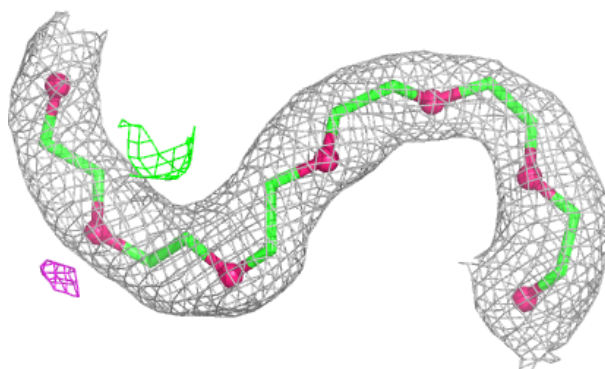
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around P6G B 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around P6G C 304:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.