



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:06 PM UTC

PDB ID : 3EUQ / pdb_00003euq
Title : X-ray structural of a type III pentaketide synthase from *Neurospora crassa*
Authors : Zhang, H.; Brunzelle, J.S.; Nair, S.K.
Deposited on : 2008-10-10
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

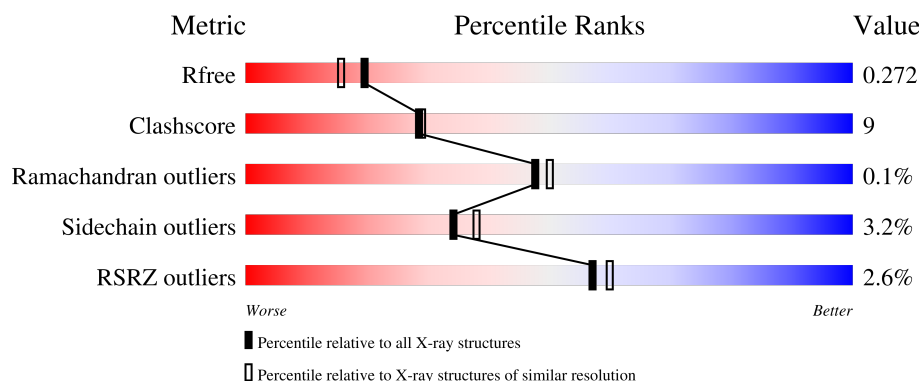
1 Overall quality at a glance

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	379	80% 18% .
1	B	379	3% 85% 14% .
1	C	379	3% 86% 13% .
1	D	379	3% 83% 16% .

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
1	CSD	A	152	-	-	X	-
1	CSD	B	152	-	-	X	-
1	CSD	C	152	-	-	X	-
1	CSD	D	152	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12262 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.

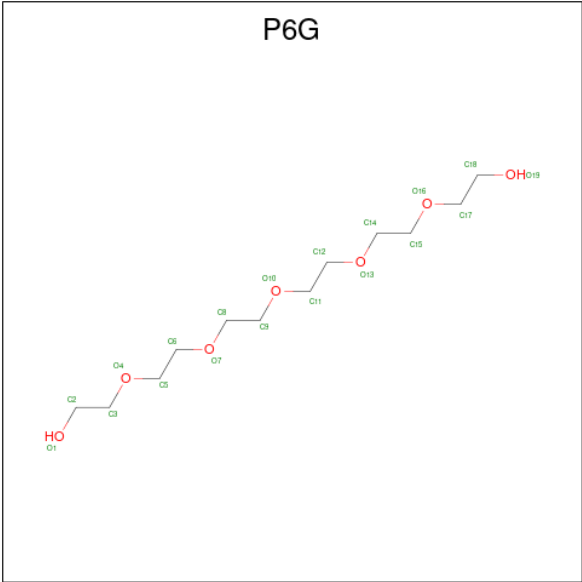
- Molecule 1 is a protein called Putative uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2850	1791	498	543	18			
1	B	379	Total	C	N	O	S	0	0	0
			2850	1791	498	543	18			
1	C	379	Total	C	N	O	S	0	0	0
			2850	1791	498	543	18			
1	D	379	Total	C	N	O	S	0	0	0
			2850	1791	498	543	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	252	GLY	PHE	engineered mutation	UNP Q7S6N4
B	252	GLY	PHE	engineered mutation	UNP Q7S6N4
C	252	GLY	PHE	engineered mutation	UNP Q7S6N4
D	252	GLY	PHE	engineered mutation	UNP Q7S6N4

- Molecule 2 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: C₁₂H₂₆O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			19	12	7		
2	C	1	Total	C	O	0	0
			19	12	7		

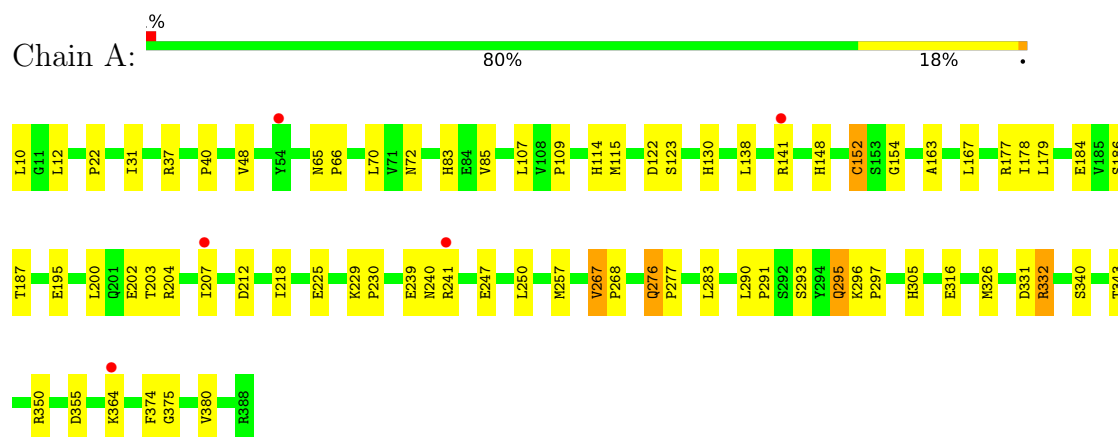
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	219	Total	O	0	0
			219	219		
3	B	203	Total	O	0	0
			203	203		
3	C	190	Total	O	0	0
			190	190		
3	D	212	Total	O	0	0
			212	212		

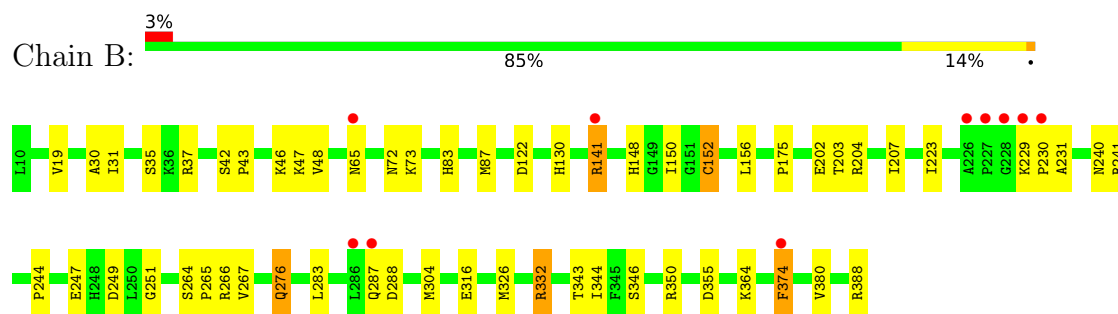
3 Residue-property plots [i](#)

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.

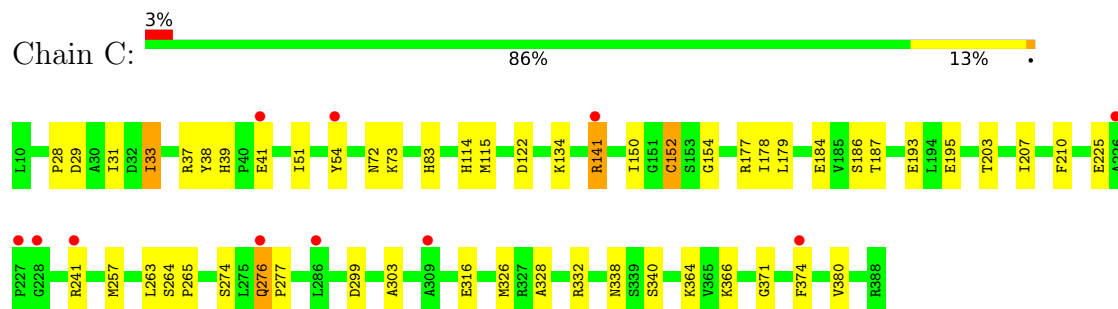
- Molecule 1: Putative uncharacterized protein



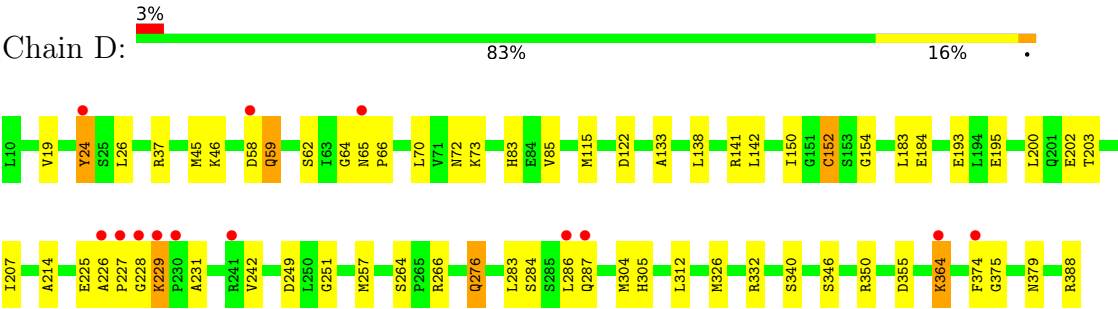
- Molecule 1: Putative uncharacterized protein



- Molecule 1: Putative uncharacterized protein



- Molecule 1: Putative uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.69Å 105.12Å 105.08Å 90.00° 90.35° 90.00°	Depositor
Resolution (Å)	25.00 – 2.10 25.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.3 (25.00-2.10) 96.2 (25.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.198 , 0.256 (Not available) , 0.272	Depositor DCC
R_{free} test set	4264 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.035 for -h,-l,-k 0.035 for -h,l,k 0.358 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12262	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: P6G, CSD

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	0.63	1/2902 (0.0%)	0.88	1/3944 (0.0%)
1	B	0.65	0/2902	0.90	4/3944 (0.1%)
1	C	0.64	0/2902	0.86	0/3944
1	D	0.65	0/2902	0.90	4/3944 (0.1%)
All	All	0.64	1/11608 (0.0%)	0.89	9/15776 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	267	VAL	CA-CB	7.19	1.57	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	225	GLU	N-CA-C	-7.51	98.33	110.20
1	A	154	GLY	N-CA-C	5.75	119.62	112.73
1	B	87	MET	N-CA-C	5.47	117.05	111.14
1	B	251	GLY	N-CA-C	5.46	118.22	110.74
1	B	150	ILE	N-CA-C	5.38	118.90	113.47
1	B	267	VAL	N-CA-CB	5.26	114.04	110.52
1	D	375	GLY	CA-C-N	5.23	125.23	119.90
1	D	375	GLY	C-N-CA	5.23	125.23	119.90
1	D	251	GLY	N-CA-C	5.15	118.06	110.80

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	2850	0	2864	67	0
1	B	2850	0	2864	54	0
1	C	2850	0	2864	47	0
1	D	2850	0	2864	48	0
2	A	19	0	26	2	0
2	C	19	0	26	2	0
3	A	219	0	0	10	0
3	B	203	0	0	7	0
3	C	190	0	0	8	0
3	D	212	0	0	5	0
All	All	12262	0	11508	201	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 (201) 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:C:152:CSD:OD2	1:C:152:CSD:SG	1.95	1.25
1:B:152:CSD:SG	1:B:152:CSD:OD2	1.95	1.25
1:D:152:CSD:SG	1:D:152:CSD:OD2	1.95	1.24
1:A:152:CSD:SG	1:A:152:CSD:OD2	1.95	1.22
1:B:152:CSD:SG	1:B:152:CSD:OD1	1.98	1.22
1:D:152:CSD:SG	1:D:152:CSD:OD1	1.99	1.21
1:C:152:CSD:SG	1:C:152:CSD:OD1	1.99	1.20
1:A:152:CSD:SG	1:A:152:CSD:OD1	1.99	1.19
1:A:167:LEU:HG	3:A:991:HOH:O	1.45	1.15
1:D:58:ASP:HB2	3:D:712:HOH:O	1.54	1.07
1:C:115:MET:HE3	1:C:179:LEU:HD23	1.38	1.05
1:D:115:MET:HE3	1:D:142:LEU:HD21	1.40	1.02
1:B:276:GLN:HE21	1:B:276:GLN:H	1.14	0.91
1:C:276:GLN:HE21	1:C:276:GLN:H	1.13	0.91
1:B:316:GLU:HG3	1:B:326:MET:HE1	1.53	0.91
1:A:115:MET:HE2	1:A:179:LEU:HD23	1.52	0.89
1:C:177:ARG:NH2	1:C:225:GLU:HG2	1.87	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:ASP:O	1:C:33:ILE:HD13	1.72	0.88
1:C:28:PRO:HA	3:C:782:HOH:O	1.74	0.88
1:D:364:LYS:HA	1:D:364:LYS:CE	2.06	0.86
1:A:276:GLN:H	1:A:276:GLN:HE21	1.27	0.81
1:B:152:CSD:OD2	1:B:152:CSD:N	2.16	0.78
1:B:65:ASN:HB3	3:B:703:HOH:O	1.85	0.75
1:D:364:LYS:HA	1:D:364:LYS:HE2	1.67	0.74
1:C:134:LYS:HD3	3:C:691:HOH:O	1.88	0.74
1:D:276:GLN:HE21	1:D:276:GLN:H	1.36	0.74
1:A:141:ARG:HH21	1:B:240:ASN:H	1.33	0.73
1:C:141:ARG:HH11	1:C:141:ARG:HB3	1.54	0.73
1:A:374:PHE:CD1	1:A:380:VAL:HG22	2.23	0.73
1:A:152:CSD:OD2	1:A:152:CSD:CB	2.37	0.73
1:A:240:ASN:H	1:B:141:ARG:NH2	1.87	0.73
1:B:316:GLU:HG3	1:B:326:MET:CE	2.19	0.73
1:B:83:HIS:HD2	1:B:122:ASP:OD1	1.71	0.73
1:C:115:MET:HE3	1:C:179:LEU:CD2	2.18	0.72
1:C:152:CSD:OD2	1:C:152:CSD:CB	2.37	0.72
1:D:115:MET:HE1	1:D:133:ALA:HB2	1.71	0.71
3:A:987:HOH:O	1:D:65:ASN:HB3	1.89	0.70
1:B:304:MET:SD	1:B:374:PHE:HZ	2.16	0.69
1:A:163:ALA:O	3:A:991:HOH:O	2.11	0.68
1:D:287:GLN:HA	3:D:668:HOH:O	1.93	0.67
1:D:364:LYS:HA	1:D:364:LYS:HE3	1.77	0.66
1:D:37:ARG:HH11	1:D:72:ASN:HD21	1.43	0.65
1:C:364:LYS:HE2	3:C:729:HOH:O	1.96	0.65
1:B:276:GLN:H	1:B:276:GLN:NE2	1.90	0.65
1:C:83:HIS:HD2	1:C:122:ASP:OD1	1.81	0.63
1:B:207:ILE:HG12	3:B:604:HOH:O	1.98	0.63
1:D:195:GLU:CD	1:D:257:MET:HE1	2.24	0.63
1:C:316:GLU:HG3	1:C:326:MET:HE1	1.80	0.62
1:B:264:SER:HB2	1:B:265:PRO:HD2	1.82	0.61
1:D:115:MET:CE	1:D:138:LEU:HD12	2.30	0.61
1:A:12:LEU:HD21	3:A:991:HOH:O	2.01	0.61
1:A:83:HIS:HD2	1:A:122:ASP:OD1	1.84	0.61
1:A:152:CSD:OD2	1:A:152:CSD:HB3	2.00	0.61
1:A:37:ARG:HH11	1:A:72:ASN:HD21	1.49	0.61
1:D:152:CSD:OD2	1:D:152:CSD:CB	2.48	0.60
1:C:187:THR:HA	2:C:600:P6G:H111	1.83	0.60
1:C:195:GLU:OE1	1:C:257:MET:HE1	2.01	0.60
1:C:152:CSD:OD2	1:C:152:CSD:HB3	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:LYS:HD2	1:A:230:PRO:HD2	1.83	0.59
1:B:152:CSD:OD2	1:B:152:CSD:CB	2.50	0.59
1:B:380:VAL:HG13	3:B:677:HOH:O	2.02	0.59
1:C:115:MET:CE	1:C:179:LEU:HD23	2.24	0.59
1:A:374:PHE:HD1	1:A:380:VAL:HG22	1.70	0.57
1:C:241:ARG:HG3	3:C:696:HOH:O	2.03	0.57
1:B:276:GLN:HE21	1:B:276:GLN:N	1.94	0.57
1:D:226:ALA:H	1:D:229:LYS:HE3	1.69	0.57
1:C:316:GLU:HG3	1:C:326:MET:CE	2.35	0.56
1:D:305:HIS:HB3	3:D:578:HOH:O	2.06	0.56
1:D:350:ARG:NH1	1:D:355:ASP:OD1	2.39	0.56
1:B:240:ASN:HB2	3:B:548:HOH:O	2.06	0.56
1:D:65:ASN:HB2	1:D:66:PRO:HD2	1.87	0.56
1:D:65:ASN:HB3	3:D:528:HOH:O	2.06	0.55
1:C:37:ARG:HH11	1:C:72:ASN:HD21	1.54	0.55
1:C:374:PHE:HD1	1:C:380:VAL:HG22	1.69	0.55
1:A:332:ARG:HH11	1:A:332:ARG:HA	1.72	0.55
1:A:276:GLN:HB2	1:A:277:PRO:HD3	1.88	0.55
1:D:195:GLU:OE2	1:D:257:MET:HE1	2.07	0.54
1:B:37:ARG:HH11	1:B:72:ASN:HD21	1.55	0.54
1:A:283:LEU:HD21	1:A:295:GLN:O	2.08	0.54
1:C:177:ARG:HH21	1:C:225:GLU:HG2	1.72	0.54
1:A:48:VAL:HG11	1:A:207:ILE:HD11	1.90	0.54
1:A:10:LEU:N	3:A:978:HOH:O	2.40	0.53
1:A:186:SER:HB2	2:A:500:P6G:H52	1.89	0.53
1:A:109:PRO:HB3	1:A:138:LEU:HD23	1.89	0.53
1:A:240:ASN:H	1:B:141:ARG:HH21	1.54	0.53
1:B:229:LYS:HD2	1:B:230:PRO:HD2	1.89	0.53
1:C:141:ARG:HB3	1:C:141:ARG:NH1	2.22	0.53
1:B:19:VAL:HG21	1:B:346:SER:HA	1.89	0.53
1:D:24:TYR:CE1	1:D:64:GLY:HA2	2.44	0.53
1:B:141:ARG:HH11	1:B:141:ARG:HB3	1.73	0.52
1:A:247:GLU:OE2	1:B:130:HIS:HE1	1.91	0.52
1:C:374:PHE:CD1	1:C:380:VAL:HG22	2.44	0.52
1:D:184:GLU:HG3	1:D:340:SER:HB3	1.92	0.52
1:A:241:ARG:HH21	1:B:141:ARG:HD2	1.75	0.52
1:B:244:PRO:HD2	1:C:54:TYR:CZ	2.46	0.52
1:A:202:GLU:OE2	1:A:204:ARG:HD2	2.11	0.51
1:B:374:PHE:CD2	1:B:374:PHE:N	2.78	0.51
1:B:231:ALA:O	1:B:388:ARG:HG3	2.11	0.51
1:B:350:ARG:NH1	1:B:355:ASP:OD1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:115:MET:HE1	1:D:138:LEU:HD12	1.92	0.51
1:B:43:PRO:O	1:B:47:LYS:HG2	2.11	0.51
1:B:276:GLN:OE1	3:B:705:HOH:O	2.19	0.51
1:C:186:SER:HB2	2:C:600:P6G:H52	1.92	0.51
1:A:12:LEU:HD11	3:A:991:HOH:O	2.10	0.51
1:D:83:HIS:HD2	1:D:122:ASP:OD1	1.94	0.51
1:A:239:GLU:HA	1:B:141:ARG:HH21	1.76	0.50
1:D:242:VAL:HG22	1:D:379:ASN:OD1	2.12	0.50
1:A:115:MET:CE	1:A:138:LEU:HD12	2.41	0.50
1:A:203:THR:HG23	3:A:771:HOH:O	2.11	0.50
1:B:31:ILE:O	1:B:35:SER:HB2	2.11	0.50
1:C:276:GLN:H	1:C:276:GLN:NE2	1.94	0.50
1:A:202:GLU:OE2	1:A:204:ARG:NH1	2.42	0.49
1:A:107:LEU:HD11	1:A:218:ILE:HG21	1.92	0.49
1:C:83:HIS:HE1	3:C:657:HOH:O	1.94	0.49
1:C:203:THR:HG21	3:C:662:HOH:O	2.10	0.49
1:D:115:MET:HE1	1:D:133:ALA:CB	2.41	0.49
1:C:274:SER:O	1:C:277:PRO:HD2	2.13	0.49
1:A:148:HIS:HB2	1:B:148:HIS:HB2	1.94	0.49
1:B:42:SER:O	1:B:46:LYS:HG3	2.12	0.49
1:C:264:SER:HB2	1:C:265:PRO:HD2	1.95	0.49
1:A:195:GLU:OE1	1:A:257:MET:HE1	2.12	0.48
1:A:200:LEU:HB2	1:A:202:GLU:HG2	1.95	0.48
1:A:141:ARG:HD2	1:B:241:ARG:HE	1.78	0.48
1:B:249:ASP:OD1	1:B:266:ARG:HD3	2.13	0.48
1:C:39:HIS:HE1	3:C:607:HOH:O	1.95	0.48
1:C:150:ILE:HD12	1:C:154:GLY:HA2	1.96	0.48
1:D:312:LEU:O	1:D:326:MET:HE1	2.13	0.48
1:A:123:SER:OG	1:B:148:HIS:HE1	1.97	0.47
1:B:229:LYS:HG3	1:B:230:PRO:O	2.14	0.47
1:A:10:LEU:HA	3:A:978:HOH:O	2.15	0.47
1:A:332:ARG:N	1:A:332:ARG:HD2	2.30	0.47
1:A:10:LEU:CA	3:A:978:HOH:O	2.63	0.47
1:A:331:ASP:OD2	1:A:332:ARG:NH1	2.48	0.46
1:A:187:THR:HA	2:A:500:P6G:H111	1.97	0.46
1:A:267:VAL:HB	1:A:268:PRO:HD3	1.97	0.46
1:D:45:MET:HB3	3:D:593:HOH:O	2.16	0.46
1:D:304:MET:HE3	1:D:304:MET:HB2	1.73	0.46
1:A:70:LEU:HD22	1:A:85:VAL:HG21	1.98	0.46
1:C:276:GLN:HB2	1:C:277:PRO:HD3	1.97	0.45
1:B:83:HIS:HE1	3:B:558:HOH:O	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:115:MET:CE	1:D:142:LEU:HD21	2.28	0.45
1:D:152:CSD:OD2	1:D:152:CSD:HB3	2.16	0.45
1:A:141:ARG:NH2	1:B:240:ASN:H	2.09	0.45
1:D:193:GLU:CB	1:D:207:ILE:HG23	2.46	0.45
1:C:195:GLU:CD	1:C:257:MET:HE1	2.42	0.45
1:A:31:ILE:HD12	1:A:207:ILE:HD12	1.99	0.45
1:A:10:LEU:N	3:A:970:HOH:O	2.49	0.44
1:D:70:LEU:HD22	1:D:85:VAL:HG21	1.99	0.44
1:A:130:HIS:CE1	1:B:247:GLU:CG	3.00	0.44
1:A:114:HIS:HB2	1:A:178:ILE:HG23	1.99	0.44
1:A:296:LYS:HB2	1:A:297:PRO:HD2	1.99	0.44
1:A:22:PRO:HG3	1:A:212:ASP:HB2	2.00	0.44
1:A:332:ARG:HA	1:A:332:ARG:NH1	2.32	0.44
1:D:150:ILE:HD12	1:D:154:GLY:HA2	1.99	0.44
1:D:19:VAL:HG21	1:D:346:SER:HA	2.00	0.44
1:D:24:TYR:CD2	1:D:24:TYR:N	2.85	0.44
1:D:183:LEU:HD13	1:D:214:ALA:HB2	2.00	0.44
1:A:250:LEU:HD23	1:A:375:GLY:C	2.43	0.43
1:D:59:GLN:HE21	1:D:59:GLN:HA	1.82	0.43
1:B:202:GLU:HG3	1:B:204:ARG:HG3	1.99	0.43
1:C:210:PHE:HA	1:C:338:ASN:O	2.17	0.43
1:A:184:GLU:HG3	1:A:340:SER:HB3	1.99	0.43
1:C:328:ALA:HB3	3:C:644:HOH:O	2.17	0.43
1:B:202:GLU:HG3	1:B:204:ARG:CG	2.49	0.43
1:B:332:ARG:HA	1:B:332:ARG:HD2	1.72	0.43
1:A:152:CSD:OD1	1:A:305:HIS:HE1	2.02	0.42
1:C:184:GLU:HG3	1:C:340:SER:HB3	1.99	0.42
1:C:299:ASP:O	1:C:366:LYS:HD2	2.19	0.42
1:A:316:GLU:HG3	1:A:326:MET:CE	2.49	0.42
1:A:40:PRO:HG2	1:D:24:TYR:OH	2.18	0.42
1:B:156:LEU:HD21	1:B:344:ILE:HG12	2.01	0.42
1:B:30:ALA:HB3	3:B:586:HOH:O	2.19	0.42
1:D:115:MET:CE	1:D:138:LEU:CD1	2.97	0.42
1:D:249:ASP:HA	1:D:264:SER:HB3	2.02	0.42
1:B:48:VAL:HG11	1:B:207:ILE:HD11	2.01	0.42
1:B:283:LEU:O	1:B:287:GLN:N	2.51	0.42
1:D:200:LEU:HB2	1:D:202:GLU:HG2	2.01	0.42
1:B:175:PRO:HB3	1:B:223:ILE:O	2.20	0.42
1:C:51:ILE:HA	1:C:54:TYR:CD2	2.54	0.42
1:A:239:GLU:HA	1:B:141:ARG:NH2	2.35	0.42
1:A:115:MET:CE	1:A:138:LEU:CD1	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:MET:HE2	1:B:304:MET:HB2	1.79	0.41
1:A:177:ARG:NH2	1:A:225:GLU:HG2	2.36	0.41
1:C:115:MET:CE	1:C:179:LEU:CD2	2.92	0.41
1:A:290:LEU:HA	1:A:291:PRO:HD3	1.95	0.41
1:C:303:ALA:O	1:C:371:GLY:HA2	2.19	0.41
1:A:350:ARG:NH1	1:A:355:ASP:OD1	2.53	0.41
1:C:274:SER:C	1:C:277:PRO:HD2	2.46	0.41
1:D:231:ALA:O	1:D:388:ARG:HG3	2.20	0.41
1:C:193:GLU:CB	1:C:207:ILE:HG23	2.51	0.41
1:D:24:TYR:CD1	1:D:64:GLY:HA2	2.56	0.41
1:A:65:ASN:HB2	1:A:66:PRO:CD	2.51	0.41
1:A:130:HIS:CE1	1:B:247:GLU:HG3	2.56	0.41
1:A:305:HIS:HD2	1:A:343:THR:HB	1.85	0.41
1:B:304:MET:SD	1:B:374:PHE:CZ	3.06	0.41
1:D:26:LEU:HG	1:D:62:SER:HB2	2.03	0.41
1:C:38:TYR:OH	1:C:72:ASN:HA	2.22	0.40
1:D:249:ASP:OD1	1:D:266:ARG:HD3	2.20	0.40
1:A:295:GLN:HE21	1:A:295:GLN:HB3	1.70	0.40
1:C:114:HIS:HB2	1:C:178:ILE:HG23	2.03	0.40
1:D:364:LYS:CE	1:D:364:LYS:CA	2.83	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	376/379 (99%)	367 (98%)	9 (2%)	0	100	100
1	B	376/379 (99%)	367 (98%)	9 (2%)	0	100	100
1	C	376/379 (99%)	367 (98%)	9 (2%)	0	100	100
1	D	376/379 (99%)	364 (97%)	10 (3%)	2 (0%)	24	22
All	All	1504/1516 (99%)	1465 (97%)	37 (2%)	2 (0%)	48	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	227	PRO
1	D	228	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	201/308 (65%)	196 (98%)	5 (2%)	42	48
1	B	308/308 (100%)	299 (97%)	9 (3%)	37	42
1	C	308/308 (100%)	300 (97%)	8 (3%)	40	46
1	D	308/308 (100%)	294 (96%)	14 (4%)	24	25
All	All	1125/1232 (91%)	1089 (97%)	36 (3%)	34	38

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	276	GLN
1	A	293	SER
1	A	295	GLN
1	A	332	ARG
1	A	364	LYS
1	B	73	LYS
1	B	141	ARG
1	B	203	THR
1	B	276	GLN
1	B	288	ASP
1	B	332	ARG
1	B	343	THR
1	B	364	LYS
1	B	374	PHE
1	C	31	ILE
1	C	33	ILE
1	C	41	GLU
1	C	73	LYS

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Mol	Chain	Res	Type
1	C	141	ARG
1	C	263	LEU
1	C	276	GLN
1	C	332	ARG
1	D	24	TYR
1	D	46	LYS
1	D	59	GLN
1	D	73	LYS
1	D	141	ARG
1	D	203	THR
1	D	229	LYS
1	D	276	GLN
1	D	283	LEU
1	D	284	SER
1	D	286	LEU
1	D	332	ARG
1	D	364	LYS
1	D	374	PHE

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	148	HIS
1	A	169	HIS
1	A	276	GLN
1	A	295	GLN
1	A	305	HIS
1	A	335	ASN
1	B	39	HIS
1	B	59	GLN
1	B	72	ASN
1	B	83	HIS
1	B	125	ASN
1	B	130	HIS
1	B	148	HIS
1	B	169	HIS
1	B	276	GLN
1	B	305	HIS
1	B	335	ASN
1	C	39	HIS
1	C	72	ASN
1	C	83	HIS

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Mol	Chain	Res	Type
1	C	130	HIS
1	C	148	HIS
1	C	169	HIS
1	C	240	ASN
1	C	276	GLN
1	C	289	GLN
1	C	295	GLN
1	C	305	HIS
1	D	59	GLN
1	D	72	ASN
1	D	83	HIS
1	D	169	HIS
1	D	276	GLN
1	D	295	GLN
1	D	305	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CSD	D	152	1	4,7,8	28.62	1 (25%)	1,8,10	0.80	0
1	CSD	C	152	1	4,7,8	28.44	1 (25%)	1,8,10	3.81	1 (100%)
1	CSD	A	152	1	4,7,8	28.48	1 (25%)	1,8,10	1.11	0
1	CSD	B	152	1	4,7,8	27.74	1 (25%)	1,8,10	3.80	1 (100%)

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	D	152	1	-	1/2/6/8	-
1	CSD	C	152	1	-	1/2/6/8	-
1	CSD	A	152	1	-	1/2/6/8	-
1	CSD	B	152	1	-	1/2/6/8	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	152	CSD	OD1-SG	57.22	1.99	1.47
1	A	152	CSD	OD1-SG	56.95	1.99	1.47
1	C	152	CSD	OD1-SG	56.87	1.99	1.47
1	B	152	CSD	OD1-SG	55.47	1.98	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	152	CSD	OD1-SG-CB	-3.81	98.58	105.60
1	B	152	CSD	OD1-SG-CB	-3.80	98.61	105.60

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	152	CSD	CA-CB-SG-OD1
1	B	152	CSD	CA-CB-SG-OD1
1	C	152	CSD	CA-CB-SG-OD1
1	D	152	CSD	CA-CB-SG-OD1

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	152	CSD	4	0
1	C	152	CSD	4	0
1	A	152	CSD	5	0
1	B	152	CSD	4	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	P6G	C	600	-	18,18,18	0.50	0	17,17,17	0.25	0
2	P6G	A	500	-	18,18,18	0.52	0	17,17,17	0.26	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
2	P6G	C	600	-	-	6/16/16/16	-
2	P6G	A	500	-	-	10/16/16/16	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	P6G	O7-C8-C9-O10
2	A	500	P6G	O10-C11-C12-O13
2	C	600	P6G	O16-C17-C18-O19
2	A	500	P6G	O16-C17-C18-O19
2	C	600	P6G	O1-C2-C3-O4
2	C	600	P6G	O7-C8-C9-O10

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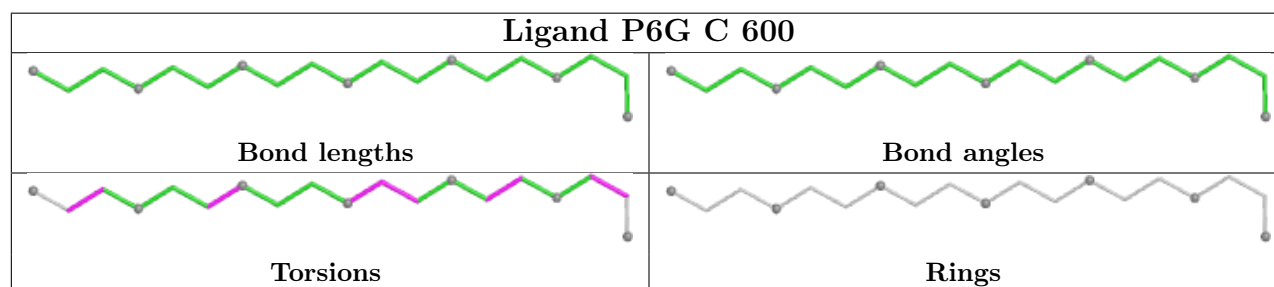
Mol	Chain	Res	Type	Atoms
2	C	600	P6G	O4-C5-C6-O7
2	A	500	P6G	C15-C14-O13-C12
2	C	600	P6G	C8-C9-O10-C11
2	C	600	P6G	C15-C14-O13-C12
2	A	500	P6G	C9-C8-O7-C6
2	A	500	P6G	C5-C6-O7-C8
2	A	500	P6G	C6-C5-O4-C3
2	A	500	P6G	C14-C15-O16-C17
2	A	500	P6G	O4-C5-C6-O7
2	A	500	P6G	C8-C9-O10-C11

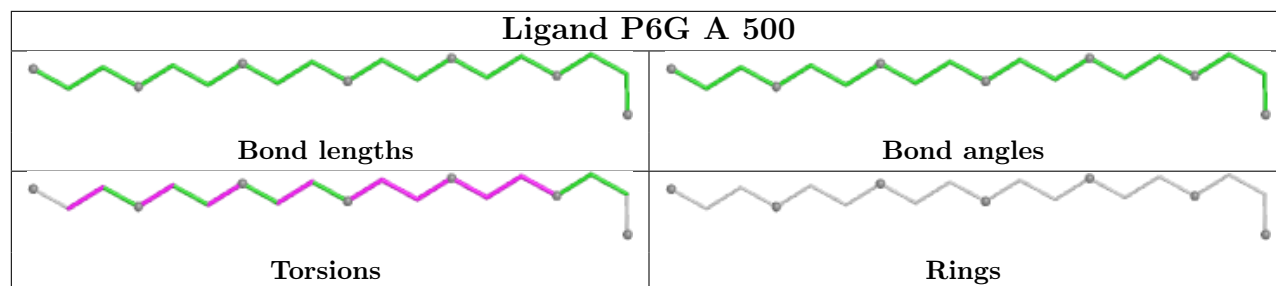
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	600	P6G	2	0
2	A	500	P6G	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	378/379 (99%)	0.22	5 (1%) 75 77	21, 27, 34, 40	0
1	B	378/379 (99%)	0.32	10 (2%) 57 60	20, 27, 35, 49	0
1	C	378/379 (99%)	0.33	11 (2%) 53 56	20, 27, 34, 42	0
1	D	378/379 (99%)	0.23	13 (3%) 48 50	20, 27, 34, 46	0
All	All	1512/1516 (99%)	0.28	39 (2%) 57 60	20, 27, 34, 49	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	227	PRO	6.8
1	D	226	ALA	5.5
1	D	227	PRO	5.3
1	D	24	TYR	4.9
1	B	226	ALA	4.7
1	C	54	TYR	4.3
1	B	287	GLN	4.2
1	B	230	PRO	4.1
1	B	374	PHE	3.8
1	D	374	PHE	3.4
1	D	228	GLY	3.2
1	C	228	GLY	3.2
1	B	229	LYS	2.9
1	D	229	LYS	2.8
1	C	141	ARG	2.8
1	D	287	GLN	2.7
1	A	241	ARG	2.6
1	B	141	ARG	2.6
1	C	309	ALA	2.5
1	C	227	PRO	2.5
1	D	241	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	228	GLY	2.5
1	A	141	ARG	2.4
1	B	65	ASN	2.3
1	D	58	ASP	2.3
1	D	65	ASN	2.3
1	A	54	TYR	2.3
1	C	286	LEU	2.3
1	D	230	PRO	2.3
1	D	286	LEU	2.3
1	D	364	LYS	2.2
1	C	41	GLU	2.2
1	C	226	ALA	2.1
1	A	207	ILE	2.1
1	A	364	LYS	2.1
1	C	374	PHE	2.0
1	B	286	LEU	2.0
1	C	276	GLN	2.0
1	C	241	ARG	2.0

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	D	152	8/9	0.89	0.11	27,29,32,34	0
1	CSD	A	152	8/9	0.93	0.09	28,30,33,35	0
1	CSD	C	152	8/9	0.94	0.09	28,29,32,37	0
1	CSD	B	152	8/9	0.94	0.09	26,27,37,37	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,

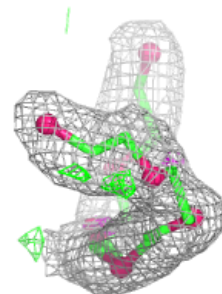
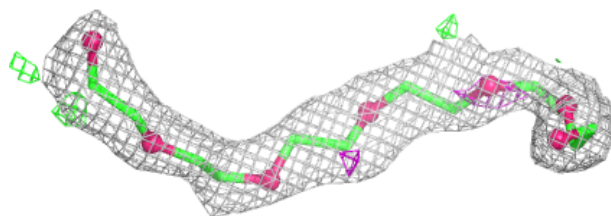
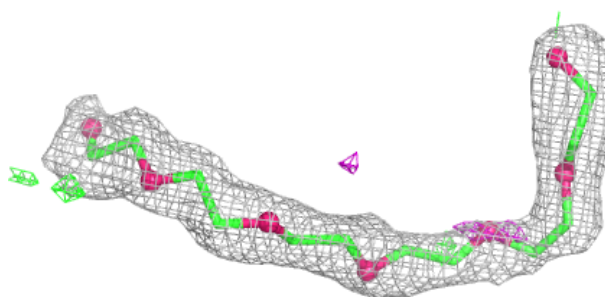
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	P6G	C	600	19/19	0.81	0.13	39,41,45,45	0
2	P6G	A	500	19/19	0.82	0.13	40,42,45,46	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.

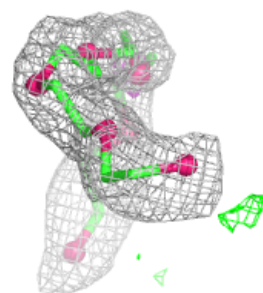
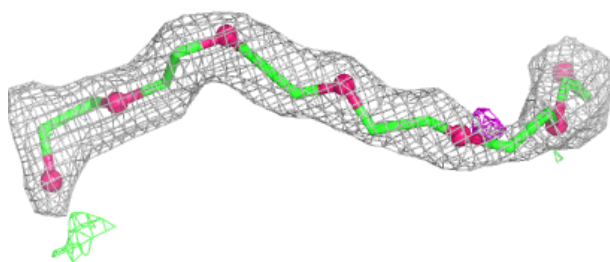
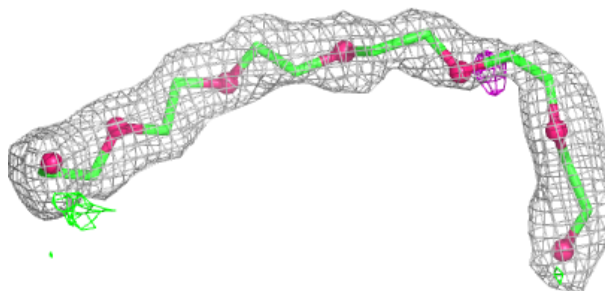
Electron density around P6G C 600:

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 A 500:

$2mF_o-DF_c$ (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.