



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:13 PM UTC

PDB ID : 4DOQ / pdb_00004doq
Title : Crystal structure of the complex of Porcine Pancreatic Trypsin with 1/2SLPI
Authors : Fukushima, K.; Takimoto-Kamimura, M.
Deposited on : 2012-02-10
Resolution : 2.00 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

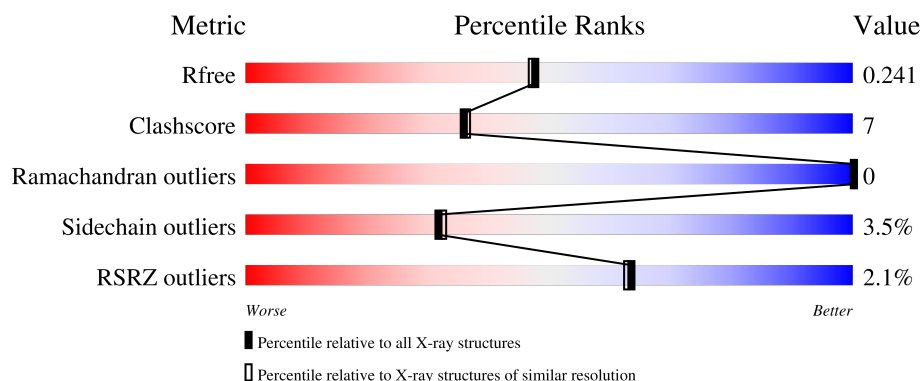
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	223	0% 83% 15% ..
1	C	223	2% 81% 16% ..
1	E	223	3% 83% 15% .
2	B	47	83% 15% .
2	D	47	2% 81% 17% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6178 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.

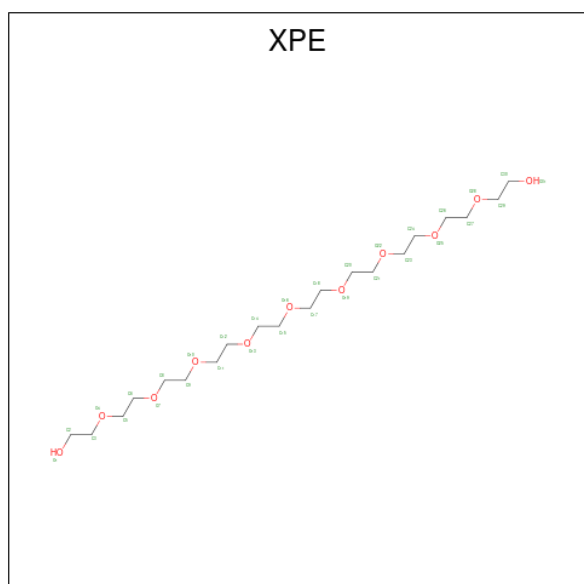
- Molecule 1 is a protein called Trypsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1630	1014	285	317	14			
1	C	219	Total	C	N	O	S	0	0	0
			1620	1009	283	314	14			
1	E	220	Total	C	N	O	S	0	0	0
			1626	1012	284	316	14			

- Molecule 2 is a protein called Antileukoproteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	47	Total	C	N	O	S	0	1	0
			352	217	61	62	12			
2	D	47	Total	C	N	O	S	0	0	0
			346	213	60	61	12			

- Molecule 3 is 3,6,9,12,15,18,21,24,27-NONAOXANONACOSANE-1,29-DIOL (CCD ID: XPE) (formula: C₂₀H₄₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			31	20	11		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

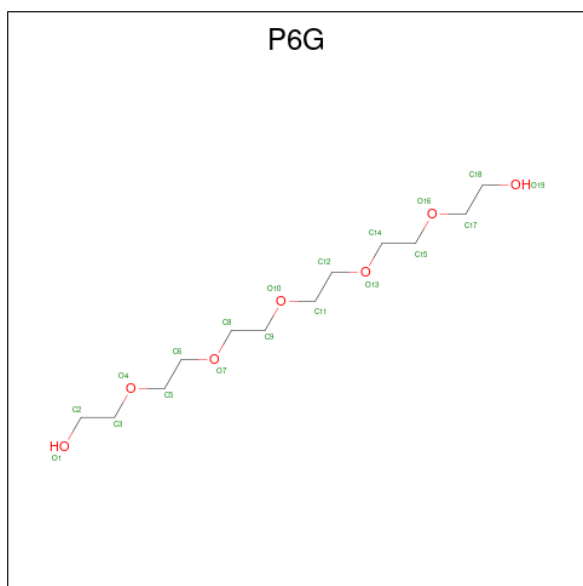
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		
5	C	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	1	Total	Ca	0	0
			1	1		

- Molecule 6 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: $C_{12}H_{26}O_7$).



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

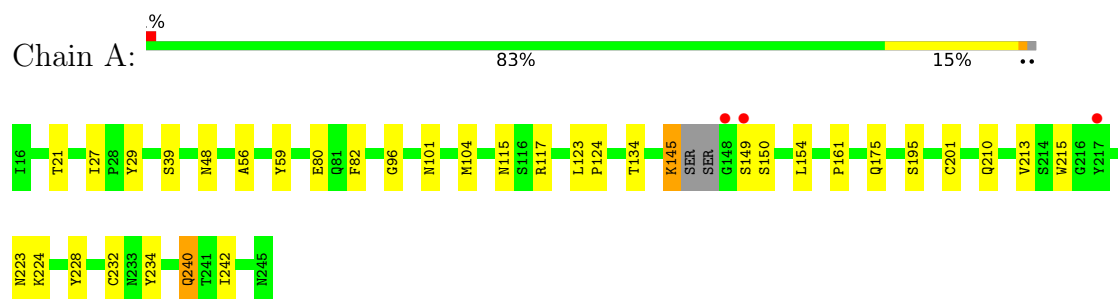
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	140	Total	O	0	0
			140	140		
7	B	41	Total	O	0	0
			41	41		
7	C	160	Total	O	0	0
			160	160		
7	D	40	Total	O	0	0
			40	40		
7	E	135	Total	O	0	0
			135	135		

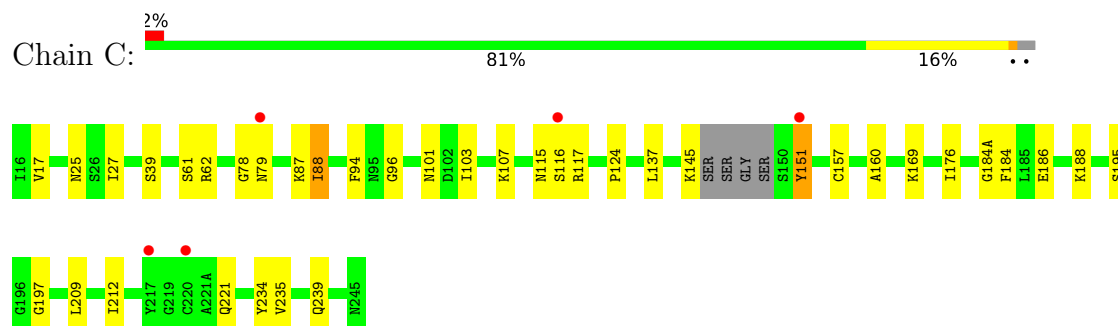
3 Residue-property plots [i](#)

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.

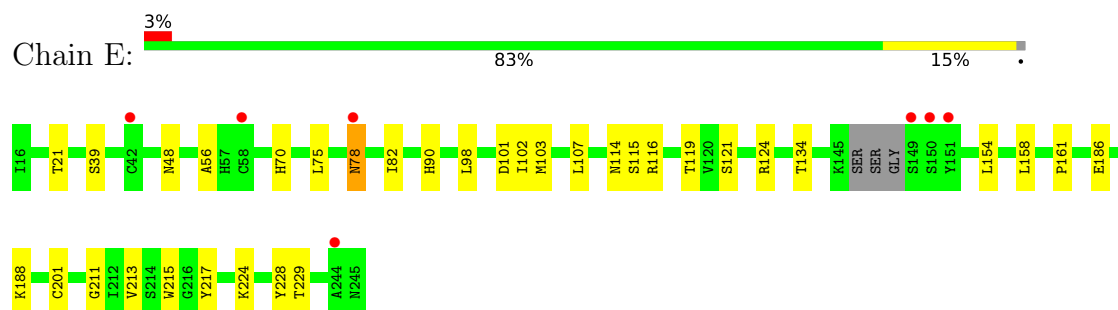
• Molecule 1: Trypsin



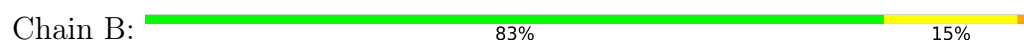
• Molecule 1: Trypsin



• Molecule 1: Trypsin

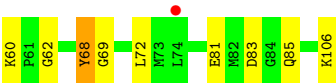
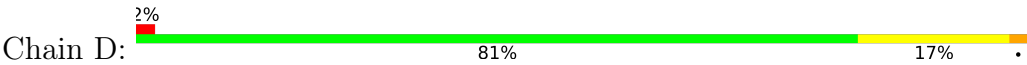


• Molecule 2: Antileukoproteinase





● Molecule 2: Antileukoproteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.53Å 118.61Å 93.39Å 90.00° 90.74° 90.00°	Depositor
Resolution (Å)	50.06 – 2.00 50.06 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.8 (50.06-2.00) 96.8 (50.06-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.187 , 0.242 0.187 , 0.241	Depositor DCC
R_{free} test set	3004 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	29.9	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6178	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, XPE, P6G, SO4

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	1.01	0/1661	1.01	4/2254 (0.2%)
1	C	1.10	0/1651	1.06	4/2241 (0.2%)
1	E	1.02	0/1657	0.98	0/2249
2	B	0.95	1/361 (0.3%)	1.29	3/481 (0.6%)
2	D	1.01	0/352	1.19	3/469 (0.6%)
All	All	1.04	1/5682 (0.0%)	1.05	14/7694 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	66	VAL	CA-CB	5.29	1.60	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	68	TYR	CA-C-N	-11.45	108.32	121.83
2	B	68	TYR	C-N-CA	-11.45	108.32	121.83
1	A	123	LEU	CA-C-N	-7.09	112.69	119.85
1	A	123	LEU	C-N-CA	-7.09	112.69	119.85
1	C	61	SER	N-CA-C	6.90	118.49	110.97
1	C	78	GLY	N-CA-C	6.40	120.43	113.58
1	C	197	GLY	CA-C-N	6.18	126.69	120.14
1	C	197	GLY	C-N-CA	6.18	126.69	120.14
2	D	69	GLY	N-CA-C	-5.86	102.20	110.75
2	B	68	TYR	O-C-N	-5.78	115.43	122.25
2	D	68	TYR	N-CA-C	5.75	118.28	110.06
2	D	62	GLY	N-CA-C	-5.74	106.19	112.04
1	A	240	GLN	CB-CA-C	-5.21	101.82	110.68
1	A	242	ILE	N-CA-C	5.10	115.32	110.42

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	1630	0	1584	30	0
1	C	1620	0	1576	23	0
1	E	1626	0	1581	17	0
2	B	352	0	353	7	0
2	D	346	0	345	5	0
3	A	31	0	42	8	0
4	A	10	0	0	0	0
4	B	5	0	0	1	0
4	C	5	0	0	0	0
4	D	10	0	0	0	0
4	E	5	0	0	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
5	E	1	0	0	0	0
6	C	19	0	26	3	0
7	A	140	0	0	4	0
7	B	41	0	0	0	0
7	C	160	0	0	4	0
7	D	40	0	0	0	0
7	E	135	0	0	1	0
All	All	6178	0	5507	81	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 7.

All (81) 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:116:SER:HB3	7:C:468:HOH:O	1.51	1.10
1:C:235:VAL:O	1:C:239:GLN:HG3	1.66	0.95
1:A:223:ASN:O	1:A:224:LYS:HE2	1.66	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:GLY:HA3	3:A:301:XPE:H172	1.56	0.85
2:D:81:GLU:H	2:D:85:GLN:NE2	1.75	0.83
1:A:149:SER:HB3	1:A:150:SER:HA	1.63	0.79
1:C:88:ILE:HD11	7:C:454:HOH:O	1.87	0.74
1:A:224:LYS:HD2	2:B:68:TYR:OH	1.89	0.72
2:D:81:GLU:H	2:D:85:GLN:HE22	1.35	0.71
2:D:68:TYR:CD2	2:D:68:TYR:O	2.43	0.71
1:C:151:TYR:N	1:C:151:TYR:CD2	2.63	0.66
1:C:94:PHE:HE2	6:C:301:P6G:H91	1.62	0.64
3:A:301:XPE:H291	7:A:455:HOH:O	2.01	0.61
3:A:301:XPE:H201	2:B:79:PHE:HE1	1.66	0.61
1:A:124:PRO:HG2	1:A:232:CYS:HA	1.84	0.59
1:A:145:LYS:NZ	1:A:145:LYS:HB2	2.18	0.58
1:E:186:GLU:H	1:E:186:GLU:CD	2.12	0.57
1:C:96:GLY:HA3	6:C:301:P6G:H122	1.87	0.55
2:B:105:VAL:HB	2:B:106:LYS:HE2	1.88	0.55
3:A:301:XPE:H171	7:A:455:HOH:O	2.07	0.55
1:A:145:LYS:NZ	1:A:149:SER:O	2.40	0.55
1:E:134:THR:O	1:E:161:PRO:HA	2.07	0.54
1:C:151:TYR:H	1:C:151:TYR:HD2	1.51	0.54
1:C:124:PRO:HG3	1:C:209:LEU:O	2.08	0.54
1:A:224:LYS:HD2	2:B:68:TYR:HH	1.73	0.53
1:A:96:GLY:HA3	3:A:301:XPE:C17	2.34	0.52
1:C:184:PHE:CE2	1:C:188:LYS:HD2	2.45	0.52
3:A:301:XPE:H202	7:A:455:HOH:O	2.10	0.51
1:E:217:TYR:HB3	1:E:224:LYS:HD2	1.93	0.51
1:A:213:VAL:HG22	1:A:228:TYR:HE2	1.75	0.50
1:C:87:LYS:HB2	1:C:107:LYS:HB3	1.93	0.50
1:C:151:TYR:N	1:C:151:TYR:HD2	2.05	0.50
1:C:62:ARG:NH1	1:E:21:THR:O	2.44	0.50
1:C:27:ILE:CD1	1:C:137:LEU:HD21	2.41	0.50
1:A:48:ASN:OD1	1:A:48:ASN:C	2.55	0.49
2:D:60:LYS:HG3	2:D:83:ASP:HB3	1.94	0.49
1:A:215:TRP:CE3	2:B:69:GLY:HA3	2.48	0.49
1:E:78:ASN:HB3	1:E:116:ARG:HD2	1.95	0.49
3:A:301:XPE:O4	3:A:301:XPE:H82	2.12	0.48
6:C:301:P6G:H52	7:C:471:HOH:O	2.13	0.48
1:E:213:VAL:HA	1:E:228:TYR:CD2	2.49	0.48
1:A:96:GLY:CA	3:A:301:XPE:H172	2.36	0.47
1:C:195:SER:OG	2:D:72:LEU:C	2.58	0.47
1:E:124:ARG:NH1	7:E:401:HOH:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:GLU:HB3	1:A:82:PHE:CE2	2.49	0.47
1:C:25:ASN:CG	1:C:117:ARG:HG3	2.40	0.46
1:A:134:THR:O	1:A:161:PRO:HA	2.15	0.46
1:E:48:ASN:HA	1:E:119:THR:HG21	1.97	0.46
1:A:101:ASN:HA	1:A:234:TYR:OH	2.16	0.46
1:A:149:SER:HB3	1:A:150:SER:CA	2.39	0.46
1:A:101:ASN:ND2	1:A:234:TYR:OH	2.48	0.45
1:E:158:LEU:HD11	1:E:188:LYS:HB3	1.99	0.45
1:C:235:VAL:O	1:C:239:GLN:CG	2.50	0.45
1:E:90:HIS:HB2	1:E:102:ILE:HG23	1.99	0.45
1:C:17:VAL:O	1:C:188:LYS:HA	2.17	0.44
1:E:82:ILE:HG21	1:E:107:LEU:HB3	2.00	0.44
1:C:103:ILE:HG13	1:C:212:ILE:CD1	2.48	0.43
1:E:213:VAL:HG22	1:E:228:TYR:HE2	1.83	0.43
1:C:27:ILE:HD11	1:C:157:CYS:SG	2.58	0.43
1:A:59:TYR:C	1:A:59:TYR:CD2	2.96	0.43
1:C:160:ALA:HB1	1:C:184(A):GLY:HA2	2.01	0.43
1:A:145:LYS:HB2	1:A:145:LYS:HZ2	1.82	0.43
1:C:87:LYS:NZ	7:C:545:HOH:O	2.49	0.43
2:B:68:TYR:HB2	4:B:201:SO4:O4	2.19	0.43
1:C:169:LYS:HG2	1:C:176:ILE:HB	2.00	0.43
1:A:27:ILE:HG23	1:A:29:TYR:CZ	2.54	0.43
1:A:39:SER:HB2	7:A:423:HOH:O	2.19	0.43
1:A:115:ASN:ND2	1:A:117:ARG:HB2	2.34	0.42
1:E:211:GLY:HA2	1:E:229:THR:O	2.19	0.42
1:A:115:ASN:HD21	1:A:117:ARG:HB2	1.84	0.42
1:C:101:ASN:HA	1:C:234:TYR:OH	2.19	0.42
1:A:21:THR:HG21	1:A:154:LEU:HD12	2.01	0.42
1:A:101:ASN:HA	1:A:234:TYR:HH	1.84	0.41
1:A:149:SER:CB	1:A:150:SER:HA	2.43	0.41
1:E:98:LEU:HD12	1:E:215:TRP:HB3	2.02	0.41
1:E:98:LEU:O	1:E:101:ASP:HB2	2.21	0.41
1:A:175:GLN:HE22	2:B:68:TYR:HD2	1.69	0.41
1:E:70:HIS:O	1:E:154:LEU:HA	2.21	0.41
1:A:124:PRO:HB3	1:A:210:GLN:NE2	2.36	0.40
1:A:56:ALA:HA	1:A:104:MET:HB3	2.04	0.40
1:E:56:ALA:HA	1:E:103:MET:HB2	2.02	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	217/223 (97%)	209 (96%)	8 (4%)	0	100	100
1	C	215/223 (96%)	205 (95%)	10 (5%)	0	100	100
1	E	216/223 (97%)	210 (97%)	6 (3%)	0	100	100
2	B	46/47 (98%)	42 (91%)	4 (9%)	0	100	100
2	D	45/47 (96%)	41 (91%)	4 (9%)	0	100	100
All	All	739/763 (97%)	707 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	181/183 (99%)	177 (98%)	4 (2%)	45	50
1	C	180/183 (98%)	172 (96%)	8 (4%)	25	24
1	E	181/183 (99%)	174 (96%)	7 (4%)	28	28
2	B	43/42 (102%)	41 (95%)	2 (5%)	23	22
2	D	42/42 (100%)	41 (98%)	1 (2%)	43	47
All	All	627/633 (99%)	605 (96%)	22 (4%)	32	32

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LYS
1	A	195	SER
1	A	201	CYS
1	A	240	GLN
2	B	73	MET
2	B	90	LEU
1	C	39	SER
1	C	79	ASN
1	C	88	ILE
1	C	115	ASN
1	C	145	LYS
1	C	151	TYR
1	C	186	GLU
1	C	221	GLN
2	D	106	LYS
1	E	39	SER
1	E	75	LEU
1	E	78	ASN
1	E	114	ASN
1	E	115	SER
1	E	121	SER
1	E	201	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	84	ASN
1	A	101	ASN
1	A	143	ASN
1	A	202	ASN
1	A	221	GLN
1	A	236	ASN
1	A	245	ASN
2	B	75	ASN
1	C	115	ASN
1	C	236	ASN
2	D	85	GLN
1	E	83	ASN
1	E	92	ASN
1	E	114	ASN
1	E	175	GLN
1	E	192	GLN

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Mol	Chain	Res	Type
1	E	236	ASN

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](#)

Of 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	201	-	4,4,4	0.32	0	6,6,6	0.28	0
4	SO4	E	301	-	4,4,4	0.33	0	6,6,6	0.18	0
4	SO4	D	201	-	4,4,4	0.23	0	6,6,6	0.33	0
4	SO4	A	303	-	4,4,4	0.21	0	6,6,6	1.08	0
4	SO4	D	202	-	4,4,4	0.25	0	6,6,6	0.40	0
4	SO4	C	302	-	4,4,4	0.29	0	6,6,6	0.10	0
6	P6G	C	301	-	18,18,18	0.47	0	17,17,17	0.65	0
3	XPE	A	301	-	30,30,30	0.52	0	29,29,29	0.46	0
4	SO4	A	302	-	4,4,4	0.41	0	6,6,6	0.18	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	XPE	A	301	-	-	15/28/28/28	-
6	P6G	C	301	-	-	9/16/16/16	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	XPE	O28-C29-C30-O31
6	C	301	P6G	O10-C11-C12-O13
3	A	301	XPE	C20-C21-O22-C23
6	C	301	P6G	O1-C2-C3-O4
6	C	301	P6G	O4-C5-C6-O7
3	A	301	XPE	C6-C5-O4-C3
3	A	301	XPE	C15-C14-O13-C12
6	C	301	P6G	C8-C9-O10-C11
6	C	301	P6G	C6-C5-O4-C3
6	C	301	P6G	C15-C14-O13-C12
3	A	301	XPE	C18-C17-O16-C15
3	A	301	XPE	C2-C3-O4-C5
3	A	301	XPE	C30-C29-O28-C27
6	C	301	P6G	C2-C3-O4-C5
6	C	301	P6G	C5-C6-O7-C8
3	A	301	XPE	O10-C11-C12-O13
3	A	301	XPE	C5-C6-O7-C8
3	A	301	XPE	O22-C23-C24-O25
3	A	301	XPE	C11-C12-O13-C14
3	A	301	XPE	C9-C8-O7-C6
3	A	301	XPE	O13-C14-C15-O16
3	A	301	XPE	O16-C17-C18-O19
3	A	301	XPE	O19-C20-C21-O22
6	C	301	P6G	O7-C8-C9-O10

There are no ring outliers.

3 monomers are involved in 12 short contacts:

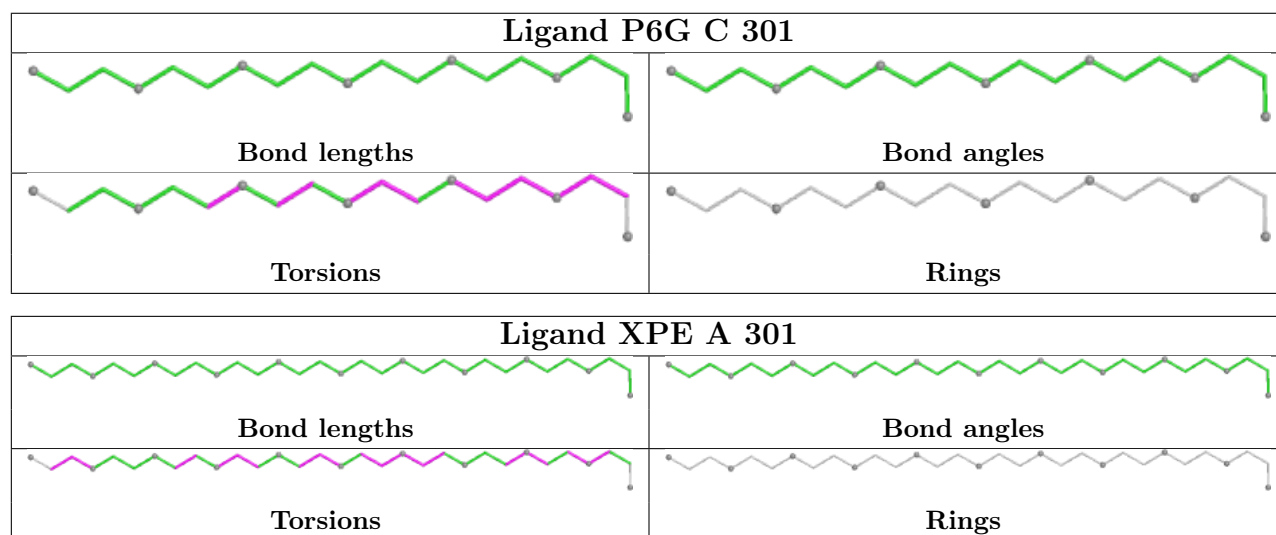
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	201	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	301	P6G	3	0
3	A	301	XPE	8	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	221/223 (99%)	0.09	3 (1%) 73 73	18, 28, 46, 63	4 (1%)
1	C	219/223 (98%)	-0.15	5 (2%) 61 60	16, 25, 42, 54	5 (2%)
1	E	220/223 (98%)	0.06	7 (3%) 50 49	10, 28, 40, 52	7 (3%)
2	B	47/47 (100%)	-0.00	0 100 100	16, 31, 40, 44	1 (2%)
2	D	47/47 (100%)	0.01	1 (2%) 63 63	23, 28, 38, 41	0
All	All	754/763 (98%)	-0.00	16 (2%) 63 63	10, 27, 42, 63	17 (2%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	149	SER	5.8
1	A	148	GLY	5.7
1	C	220	CYS	4.9
1	C	151	TYR	4.5
1	E	58	CYS	3.5
1	A	217	TYR	3.0
1	E	151	TYR	3.0
1	E	78	ASN	2.9
1	E	42	CYS	2.8
1	E	149	SER	2.6
1	C	79	ASN	2.5
1	E	150	SER	2.5
1	C	116	SER	2.4
1	C	217	TYR	2.4
1	E	244	ALA	2.3
2	D	74	LEU	2.2

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

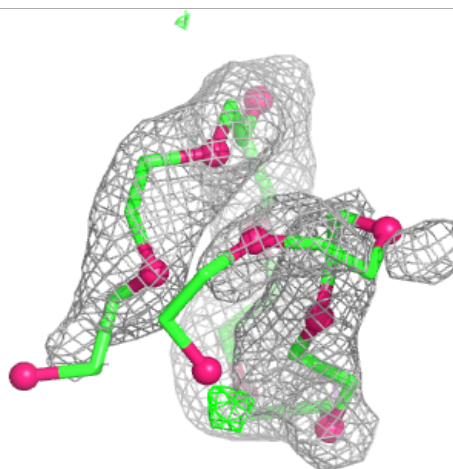
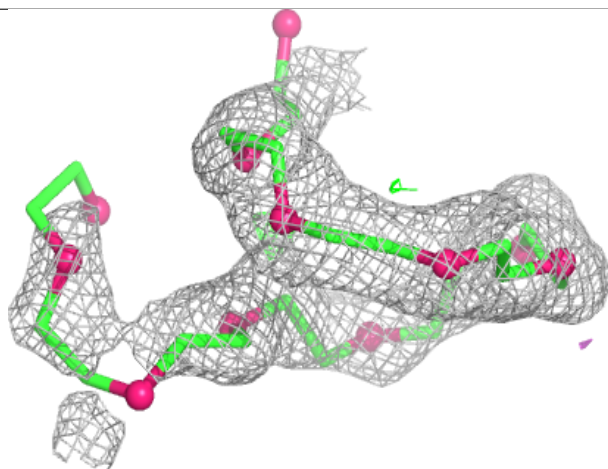
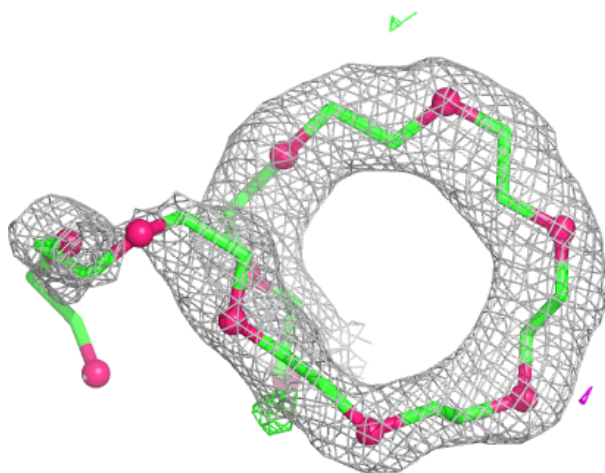
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	B	201	5/5	0.74	0.17	88,89,89,89	0
4	SO4	D	202	5/5	0.75	0.17	86,87,87,88	0
4	SO4	A	302	5/5	0.82	0.13	67,67,68,69	0
4	SO4	A	303	5/5	0.88	0.13	40,43,44,47	0
4	SO4	C	302	5/5	0.89	0.10	68,69,69,70	0
3	XPE	A	301	31/31	0.89	0.14	31,53,76,76	0
4	SO4	E	301	5/5	0.91	0.11	61,62,64,64	0
6	P6G	C	301	19/19	0.92	0.12	37,41,52,53	0
5	CA	A	304	1/1	0.96	0.06	37,37,37,37	0
4	SO4	D	201	5/5	0.97	0.07	34,37,39,41	0
5	CA	E	302	1/1	0.99	0.04	27,27,27,27	0
5	CA	C	303	1/1	0.99	0.05	37,37,37,37	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.

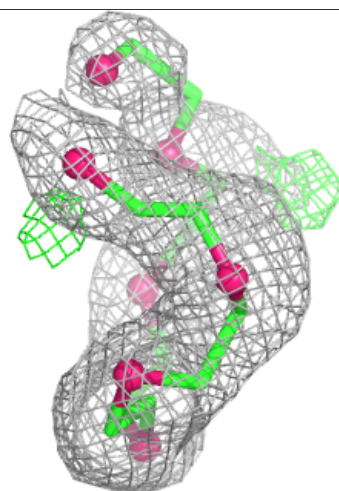
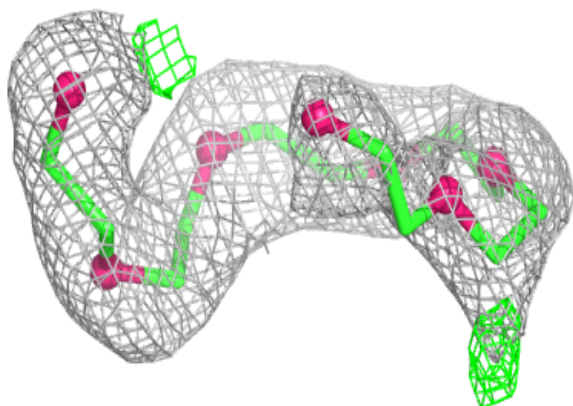
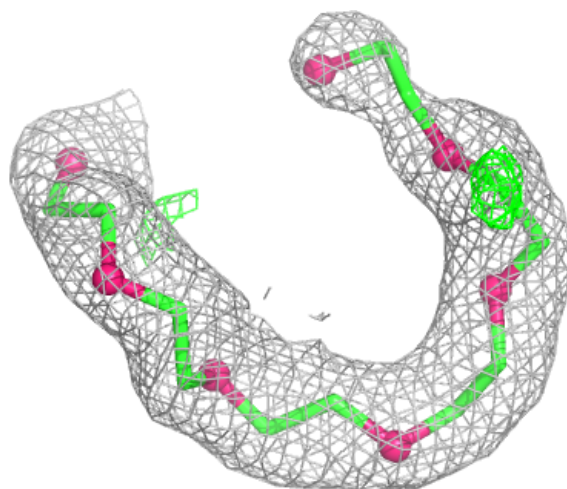
Electron density around XPE A 301:

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



Electron density around P6G C 301:

$2mF_o-DF_c$ (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.