Full wwPDB X-ray Structure Validation Report (i)

Oct 9, 2023 – 10:45 PM EDT

PDB ID	:	6XOG
Title	:	Structure of SUMO1-ML786519 adduct bound to SAE
Authors	:	Sintchak, M.; Lane, W.; Bump, N.
Deposited on	:	2020-07-07
Resolution	:	1.98 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 (i) 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 (i)) were used in the production of this report:

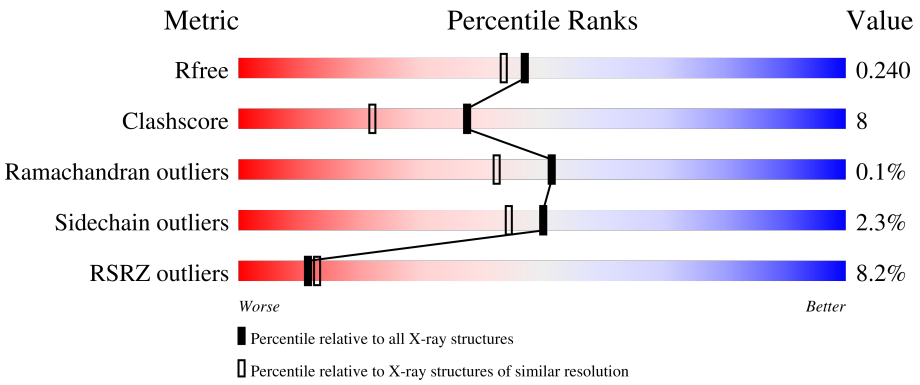
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

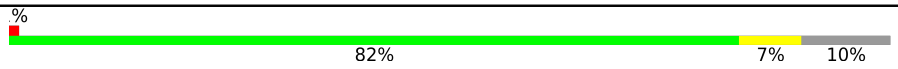
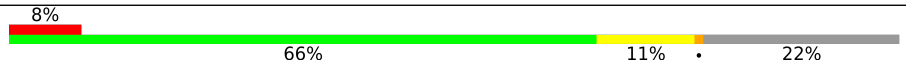
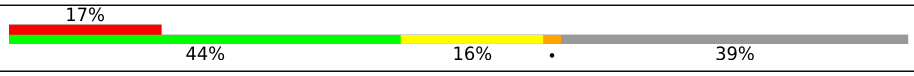
The reported resolution of this entry is 1.98 Å.

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}	130704	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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 <=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	346	
2	B	640	
3	C	101	

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 6931 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 SUMO-activating enzyme subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	311	2387	1516	407	452	12	0	0	0

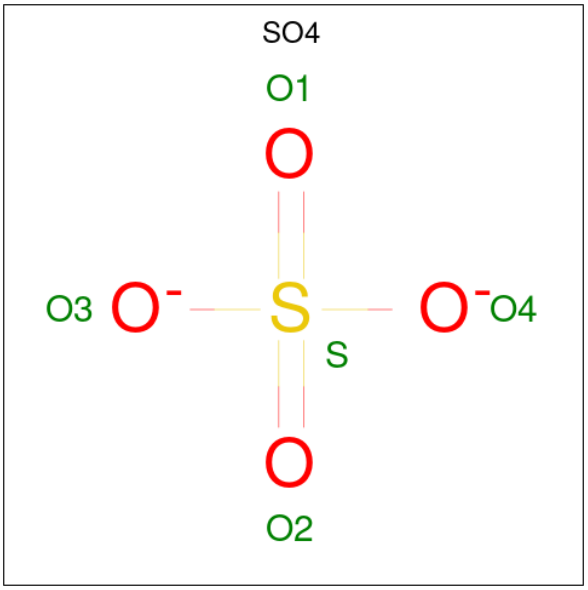
- Molecule 2 is a protein called SUMO-activating enzyme subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	500	3758	2405	632	702	19	0	0	0

- Molecule 3 is a protein called Small ubiquitin-related modifier 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	62	489	310	83	94	2	0	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

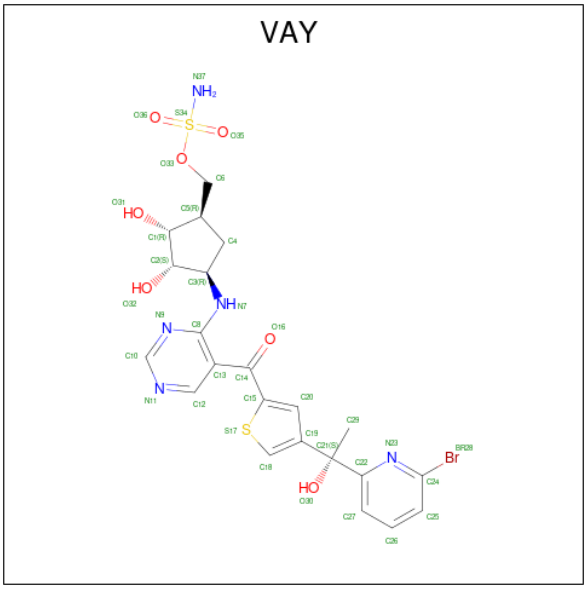


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
4	A	1	5	4	1	0	0

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
5	B	1	1	1	0	0

- Molecule 6 is {(1R,2R,3S,4R)-4-[(5-{4-[(1S)-1-(6-bromopyridin-2-yl)-1-hydroxyethyl]thiophene-2-carbonyl}pyrimidin-4-yl)amino]-2,3-dihydroxycyclopentyl}methyl sulfamate (three-letter code: VAY) (formula: $C_{22}H_{24}BrN_5O_7S_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	Br	C	N	O	S		
6	C	1	37	1	22	5	7	2	0	0

- Molecule 7 is water.

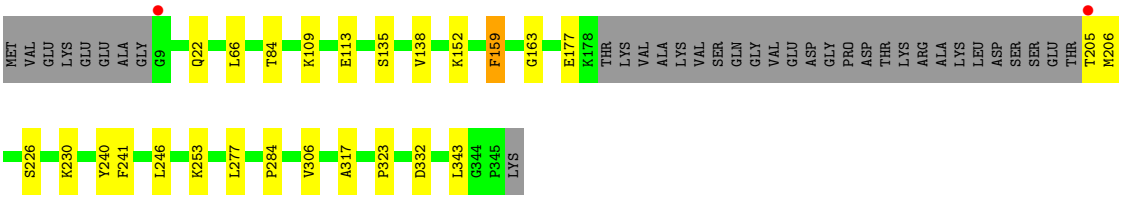
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	156	156	156	0	0
			Total	O		
7	B	86	86	86	0	0
			Total	O		
7	C	12	12	12	0	0

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.

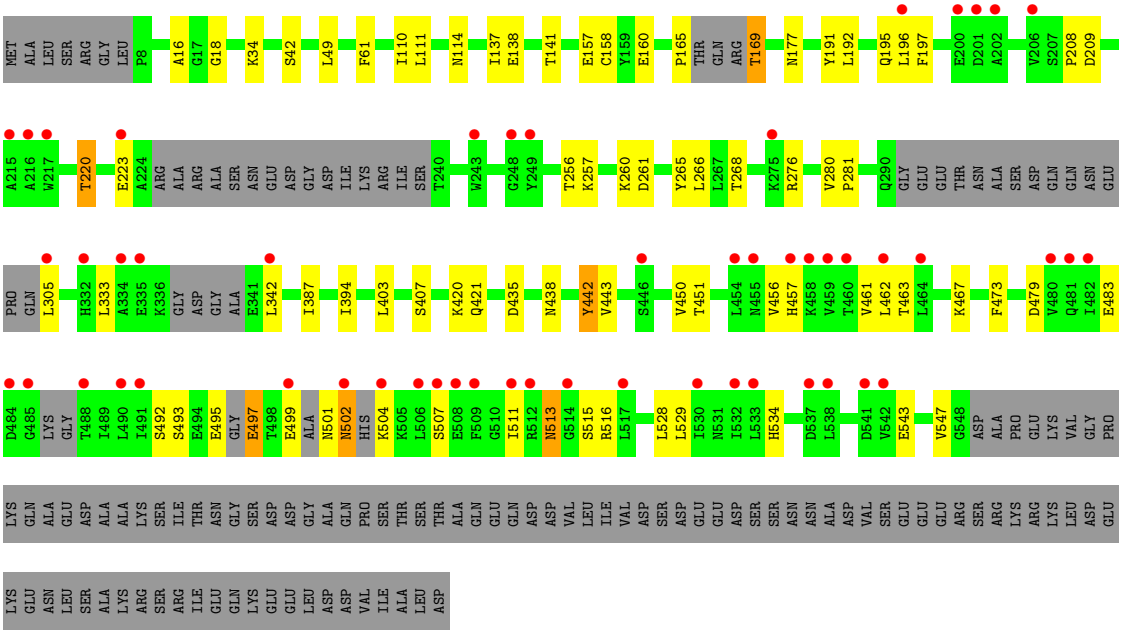
• Molecule 1: SUMO-activating enzyme subunit 1

Chain A:



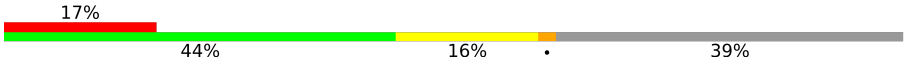
• Molecule 2: SUMO-activating enzyme subunit 2

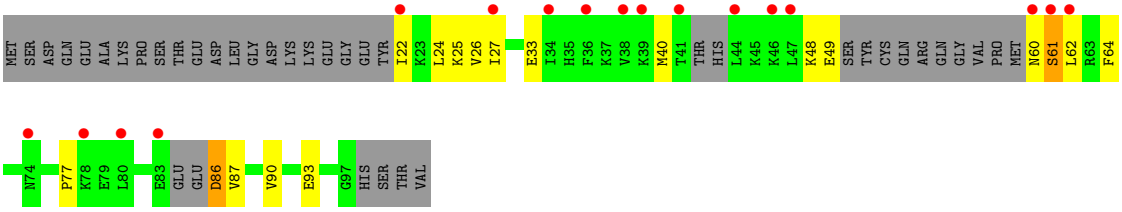
Chain B:



• Molecule 3: Small ubiquitin-related modifier 1

Chain C:




MET SER ASP GLN GLU ALA LYS PRO SER THR GLU ASP LEU GLY ASP LYS LYS GLU GLY GLU TYR I22 K23 L24 K25 V26 I27 E33 I34 H35 F36 K37 V38 K39 M40 T41 THR HIS L44 K45 K46 L47 K48 E49 SER TYR CYS GLN ARG GLN GLY VAL PRO MET N60 S61 L62 R63 F64
N74 P77 K78 E79 L80 E83 GLU GLU D86 V87 V90 E93 G97 HIS SER THR VAL

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	57.90Å 72.42Å 126.87Å 90.00° 93.86° 90.00°	Depositor
Resolution (Å)	47.70 – 1.98 47.66 – 1.98	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.70-1.98) 98.5 (47.66-1.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.36 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.206 , 0.240 0.206 , 0.240	Depositor DCC
R_{free} test set	3655 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	31.9	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 51.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6931	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: VAY, ZN, 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	0.57	1/2430 (0.0%)	0.81	1/3287 (0.0%)
2	B	0.42	1/3827 (0.0%)	0.75	1/5209 (0.0%)
3	C	0.38	0/492	0.68	0/653
All	All	0.48	2/6749 (0.0%)	0.77	2/9149 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	MET	C-N	-5.43	1.21	1.34
2	B	160	GLU	CD-OE1	5.30	1.31	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	177	GLU	CB-CA-C	5.42	121.25	110.40
2	B	420	LYS	N-CA-CB	-5.29	101.08	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	2387	0	2340	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3758	0	3654	67	0
3	C	489	0	484	26	0
4	A	5	0	0	0	0
5	B	1	0	0	0	0
6	C	37	0	0	2	0
7	A	156	0	0	5	0
7	B	86	0	0	5	0
7	C	12	0	0	1	0
All	All	6931	0	6478	101	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 8.

All (101) 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:165:PRO:HB3	3:C:61:SER:CB	1.72	1.18
3:C:93:GLU:HB2	7:C:310:HOH:O	1.49	1.11
2:B:165:PRO:HB3	3:C:61:SER:HB2	1.30	1.07
2:B:165:PRO:CB	3:C:61:SER:HB2	1.91	1.00
2:B:502:ASN:HD22	2:B:502:ASN:H	1.02	0.95
2:B:165:PRO:HB3	3:C:61:SER:HB3	1.48	0.95
2:B:499:GLU:HA	2:B:502:ASN:ND2	1.87	0.90
2:B:265:TYR:O	2:B:268:THR:HB	1.80	0.82
2:B:442:TYR:CE1	2:B:443:VAL:HG23	2.15	0.82
2:B:502:ASN:H	2:B:502:ASN:ND2	1.70	0.81
2:B:483:GLU:OE2	2:B:516:ARG:HD3	1.82	0.79
2:B:165:PRO:CG	3:C:61:SER:HB2	2.11	0.78
2:B:442:TYR:CD1	2:B:443:VAL:HG23	2.19	0.78
2:B:456:VAL:HG23	2:B:534:HIS:ND1	1.99	0.77
3:C:40:MET:HG3	3:C:77:PRO:HB2	1.69	0.75
2:B:502:ASN:HD22	2:B:502:ASN:N	1.83	0.74
2:B:165:PRO:HG3	3:C:61:SER:HB2	1.70	0.71
2:B:169:THR:HA	7:B:865:HOH:O	1.90	0.70
3:C:86:ASP:N	3:C:86:ASP:OD1	2.26	0.68
2:B:495:GLU:O	2:B:497:GLU:N	2.27	0.66
2:B:141:THR:HG22	2:B:394:ILE:HG13	1.79	0.64
2:B:499:GLU:CA	2:B:502:ASN:ND2	2.61	0.64
1:A:152:LYS:HE3	7:A:576:HOH:O	1.98	0.63
2:B:442:TYR:HE2	2:B:473:PHE:HD1	1.48	0.62
3:C:24:LEU:CD2	3:C:86:ASP:N	2.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:403:LEU:O	2:B:407:SER:HB2	2.01	0.61
1:A:205:THR:N	7:A:501:HOH:O	2.35	0.60
2:B:61:PHE:HD1	7:B:804:HOH:O	1.84	0.60
2:B:479:ASP:OD1	2:B:493:SER:N	2.36	0.59
2:B:157:GLU:OE2	2:B:435:ASP:HB2	2.01	0.59
3:C:62:LEU:HD22	3:C:90:VAL:HG12	1.85	0.58
2:B:220:THR:O	2:B:223:GLU:HG3	2.03	0.57
1:A:323:PRO:HG2	2:B:421:GLN:NE2	2.20	0.57
2:B:114:ASN:ND2	2:B:138:GLU:OE1	2.29	0.57
2:B:442:TYR:CE1	2:B:443:VAL:CG2	2.85	0.57
2:B:450:VAL:HG13	2:B:528:LEU:HD11	1.85	0.57
2:B:501:ASN:O	2:B:504:LYS:N	2.38	0.56
3:C:24:LEU:HD23	3:C:86:ASP:N	2.21	0.56
1:A:113:GLU:CB	7:A:544:HOH:O	2.54	0.56
2:B:177:ASN:HD21	2:B:209:ASP:HB3	1.70	0.56
2:B:442:TYR:HE2	2:B:473:PHE:CD1	2.24	0.56
2:B:483:GLU:OE2	2:B:516:ARG:CB	2.55	0.55
3:C:48:LYS:HB2	3:C:64:PHE:CZ	2.42	0.55
2:B:483:GLU:OE2	2:B:516:ARG:HB2	2.07	0.54
1:A:159:PHE:CG	1:A:246:LEU:HD13	2.42	0.54
2:B:333:LEU:O	2:B:333:LEU:HG	2.08	0.53
3:C:27:ILE:HD12	3:C:87:VAL:HG11	1.92	0.52
2:B:16:ALA:O	2:B:42:SER:OG	2.28	0.52
2:B:463:THR:HG21	2:B:543:GLU:OE2	2.09	0.51
1:A:163:GLY:HA3	1:A:306:VAL:HG21	1.93	0.51
2:B:461:VAL:HG12	2:B:461:VAL:O	2.10	0.51
2:B:456:VAL:HG11	2:B:511:ILE:O	2.10	0.51
2:B:49:LEU:HD23	6:C:201:VAY:C13	2.41	0.51
2:B:18:GLY:HA3	2:B:110:ILE:HD13	1.93	0.51
2:B:197:PHE:HB3	2:B:342:LEU:HD12	1.92	0.50
3:C:62:LEU:HD22	3:C:90:VAL:CG1	2.42	0.49
1:A:241:PHE:HB3	1:A:277:LEU:HD21	1.95	0.49
2:B:138:GLU:HG3	7:B:801:HOH:O	2.12	0.48
2:B:191:TYR:O	2:B:195:GLN:HG2	2.13	0.48
2:B:49:LEU:HD13	6:C:201:VAY:BR28	2.68	0.48
1:A:277:LEU:HD12	1:A:284:PRO:HA	1.95	0.47
3:C:48:LYS:HB2	3:C:64:PHE:CE1	2.49	0.46
2:B:457:HIS:ND1	2:B:507:SER:CB	2.79	0.46
2:B:499:GLU:H	2:B:502:ASN:HD21	1.64	0.46
3:C:25:LYS:HD2	3:C:33:GLU:OE2	2.16	0.46
1:A:253:LYS:HE2	7:A:563:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:457:HIS:CE1	2:B:507:SER:CB	2.99	0.46
2:B:513:ASN:HB2	2:B:534:HIS:HB3	1.98	0.46
2:B:208:PRO:HB3	2:B:266:LEU:HD13	1.97	0.45
3:C:26:VAL:HG13	3:C:90:VAL:CG2	2.47	0.45
3:C:62:LEU:HD13	3:C:90:VAL:HG11	1.99	0.45
2:B:256:THR:HG23	2:B:260:LYS:HD3	1.99	0.45
3:C:27:ILE:HD12	3:C:87:VAL:CG1	2.47	0.44
1:A:109:LYS:HD3	7:A:611:HOH:O	2.17	0.44
1:A:135:SER:OG	1:A:138:VAL:HG23	2.17	0.43
1:A:226:SER:O	1:A:230:LYS:HG3	2.18	0.43
2:B:456:VAL:H	2:B:534:HIS:CE1	2.36	0.43
2:B:483:GLU:HB2	2:B:516:ARG:O	2.19	0.43
1:A:22:GLN:HB3	2:B:387:ILE:HG22	2.01	0.42
1:A:240:TYR:CD1	1:A:343:LEU:HD13	2.54	0.42
2:B:467:LYS:NZ	7:B:809:HOH:O	2.52	0.42
2:B:492:SER:OG	2:B:497:GLU:HB3	2.20	0.42
2:B:305:LEU:HD22	7:B:874:HOH:O	2.19	0.42
2:B:257:LYS:HA	2:B:261:ASP:HB2	2.02	0.41
2:B:499:GLU:HA	2:B:502:ASN:CG	2.39	0.41
3:C:62:LEU:HB3	3:C:90:VAL:CG1	2.49	0.41
1:A:152:LYS:HE2	1:A:317:ALA:O	2.20	0.41
3:C:26:VAL:CG1	3:C:90:VAL:CG2	2.98	0.41
2:B:165:PRO:CB	3:C:61:SER:CB	2.59	0.41
2:B:499:GLU:N	2:B:502:ASN:HD21	2.19	0.41
3:C:22:ILE:HD11	3:C:77:PRO:CG	2.51	0.41
1:A:66:LEU:C	1:A:66:LEU:HD23	2.41	0.41
1:A:332:ASP:C	1:A:332:ASP:OD1	2.59	0.41
3:C:49:GLU:OE1	3:C:49:GLU:HA	2.19	0.41
2:B:192:LEU:HD11	2:B:196:LEU:HD11	2.03	0.40
2:B:451:THR:HG22	2:B:547:VAL:CG2	2.51	0.40
3:C:25:LYS:HB2	3:C:25:LYS:HE3	1.88	0.40
2:B:111:LEU:HD11	2:B:137:ILE:HG13	2.04	0.40
2:B:438:ASN:OD1	2:B:438:ASN:C	2.60	0.40
2:B:461:VAL:O	2:B:461:VAL:CG1	2.70	0.40
2:B:280:VAL:HA	2:B:281:PRO:HD3	1.95	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	307/346 (89%)	298 (97%)	9 (3%)	0	100	100
2	B	482/640 (75%)	459 (95%)	22 (5%)	1 (0%)	47	38
3	C	54/101 (54%)	51 (94%)	3 (6%)	0	100	100
All	All	843/1087 (78%)	808 (96%)	34 (4%)	1 (0%)	51	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	513	ASN

5.3.2 Protein sidechains (i)

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	254/296 (86%)	252 (99%)	2 (1%)	81	80
2	B	394/552 (71%)	383 (97%)	11 (3%)	43	32
3	C	53/91 (58%)	50 (94%)	3 (6%)	20	9
All	All	701/939 (75%)	685 (98%)	16 (2%)	50	44

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	THR
1	A	159	PHE

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Mol	Chain	Res	Type
2	B	34	LYS
2	B	158	CYS
2	B	169	THR
2	B	220	THR
2	B	276	ARG
2	B	442	TYR
2	B	462	LEU
2	B	497	GLU
2	B	502	ASN
2	B	515	SER
2	B	529	LEU
3	C	60	ASN
3	C	61	SER
3	C	86	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	195	GLN
2	B	421	GLN
2	B	502	ASN
2	B	513	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\#\|Z\| > 2$
6	VAY	C	201	3	37,40,40	2.78	10 (27%)	44,60,60	3.31	19 (43%)
4	SO4	A	401	-	4,4,4	0.30	0	6,6,6	0.13	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
6	VAY	C	201	3	-	1/20/46/46	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	201	VAY	S34-N37	9.45	1.69	1.58
6	C	201	VAY	O35-S34	6.57	1.47	1.42
6	C	201	VAY	O36-S34	6.55	1.47	1.42
6	C	201	VAY	O33-S34	-5.49	1.50	1.57
6	C	201	VAY	C3-N7	4.43	1.52	1.45
6	C	201	VAY	C15-S17	-3.26	1.68	1.72
6	C	201	VAY	O30-C21	-2.69	1.38	1.43
6	C	201	VAY	C2-C3	2.27	1.56	1.53
6	C	201	VAY	C24-N23	2.21	1.34	1.32
6	C	201	VAY	C18-S17	2.02	1.73	1.70

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	C	201	VAY	O30-C21-C19	-11.03	88.19	109.15
6	C	201	VAY	C12-C13-C8	8.24	122.61	114.55
6	C	201	VAY	O36-S34-O35	-6.08	114.21	119.97
6	C	201	VAY	C29-C21-C19	-5.10	98.24	111.34
6	C	201	VAY	N11-C10-N9	-5.04	122.15	127.70
6	C	201	VAY	C25-C24-N23	-5.03	121.51	125.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	C	201	VAY	C13-C8-N9	-4.49	115.78	121.90
6	C	201	VAY	C24-N23-C22	4.37	119.97	116.96
6	C	201	VAY	C12-N11-C10	4.05	120.98	115.80
6	C	201	VAY	C19-C18-S17	-3.89	109.55	112.25
6	C	201	VAY	N7-C8-N9	3.50	123.05	117.87
6	C	201	VAY	O31-C1-C5	-3.45	103.81	112.75
6	C	201	VAY	C4-C5-C6	-3.23	106.43	112.68
6	C	201	VAY	C26-C25-C24	3.20	119.70	116.70
6	C	201	VAY	C13-C12-N11	-3.04	117.00	123.81
6	C	201	VAY	C13-C14-C15	2.90	123.72	119.06
6	C	201	VAY	C4-C3-C2	-2.35	99.46	103.17
6	C	201	VAY	BR28-C24-C25	2.32	121.02	118.64
6	C	201	VAY	C6-O33-S34	2.07	119.80	117.21

There are no chirality outliers.

All (1) torsion outliers are listed below:

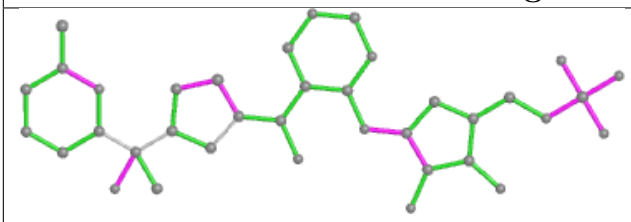
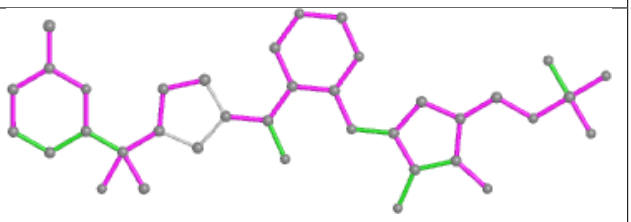
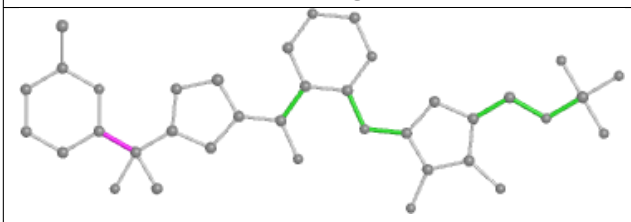
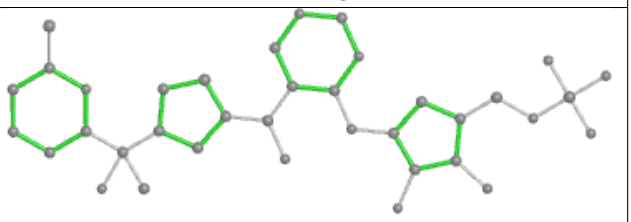
Mol	Chain	Res	Type	Atoms
6	C	201	VAY	C29-C21-C22-N23

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	201	VAY	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 then 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.

Ligand VAY C 201	
 Bond lengths	 Bond angles
 Torsions	 Rings

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	311/346 (89%)	-0.07	2 (0%)	89	90	17, 32, 58, 80	0
2	B	500/640 (78%)	0.71	53 (10%)	6	7	18, 47, 93, 114	0
3	C	62/101 (61%)	1.44	17 (27%)	0	0	23, 77, 108, 115	0
All	All	873/1087 (80%)	0.49	72 (8%)	11	13	17, 41, 92, 115	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	47	LEU	6.6
2	B	506	LEU	6.4
2	B	458	LYS	6.4
3	C	61	SER	5.9
3	C	78	LYS	5.7
2	B	532	ILE	5.4
2	B	490	LEU	5.3
3	C	22	ILE	5.0
2	B	460	THR	4.8
2	B	216	ALA	4.7
2	B	482	ILE	4.6
2	B	480	VAL	4.3
2	B	223	GLU	4.3
3	C	36	PHE	4.2
2	B	243	TRP	4.1
2	B	305	LEU	4.0
2	B	249	TYR	3.9
2	B	275	LYS	3.9
2	B	454	LEU	3.9
2	B	511	ILE	3.7
2	B	459	VAL	3.6
2	B	509	PHE	3.6
3	C	38	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
2	B	537	ASP	3.6
3	C	46	LYS	3.6
2	B	206	VAL	3.5
2	B	484	ASP	3.5
3	C	39	LYS	3.4
2	B	462	LEU	3.3
2	B	464	LEU	3.2
1	A	205	THR	3.2
2	B	517	LEU	3.2
2	B	538	LEU	3.1
2	B	202	ALA	3.1
2	B	541	ASP	3.1
2	B	481	GLN	2.9
2	B	446	SER	2.9
2	B	342	LEU	2.9
2	B	542	VAL	2.9
3	C	34	ILE	2.8
2	B	512	ARG	2.8
2	B	196	LEU	2.8
2	B	530	ILE	2.7
2	B	488	THR	2.7
3	C	80	LEU	2.7
3	C	83	GLU	2.7
2	B	502	ASN	2.6
3	C	41	THR	2.6
3	C	62	LEU	2.6
2	B	217	TRP	2.6
2	B	215	ALA	2.5
2	B	457	HIS	2.5
3	C	74	ASN	2.4
2	B	491	ILE	2.4
2	B	248	GLY	2.4
3	C	44	LEU	2.3
2	B	332	HIS	2.3
2	B	508	GLU	2.3
2	B	334	ALA	2.3
2	B	514	GLY	2.3
2	B	507	SER	2.2
2	B	485	GLY	2.2
2	B	201	ASP	2.2
2	B	455	ASN	2.1
2	B	200	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	533	LEU	2.1
2	B	499	GLU	2.1
3	C	60	ASN	2.1
1	A	9	GLY	2.0
2	B	335	GLU	2.0
2	B	504	LYS	2.0
3	C	27	ILE	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

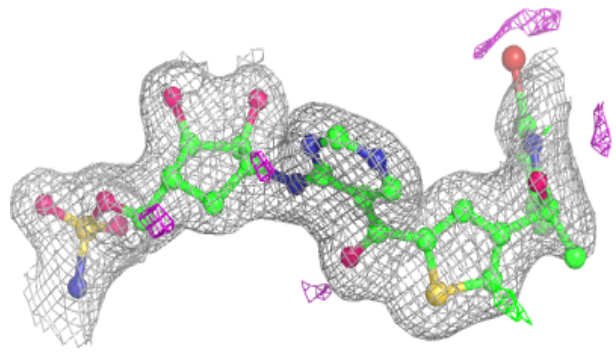
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

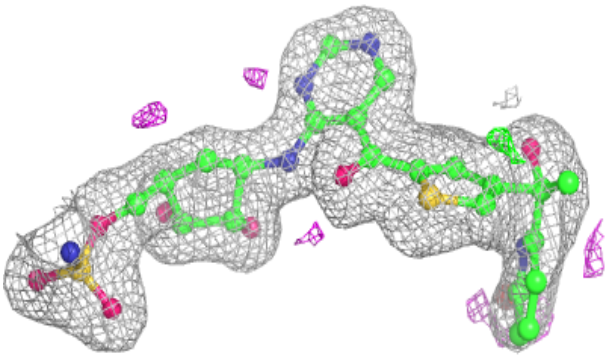
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	SO4	A	401	5/5	0.97	0.08	45,51,58,60	0
5	ZN	B	700	1/1	0.97	0.10	41,41,41,41	0
6	VAY	C	201	37/37	0.97	0.12	21,32,119,210	0

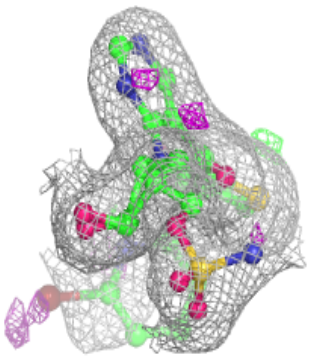
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around VAY C 201:

$2mF_o$-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.