

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110918\
 Data File : VX005886.D
 Acq On : 10 Nov 2018 03:04
 Operator : JC/MD
 Sample : J5889-12
 Misc : 20.29g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BOTTOM-C

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.666	63	84	86	rBV3	11160	41936	7.42%	0.897%
2	2.733	252	259	265	rBV2	3150	6538	1.16%	0.140%
3	3.025	299	307	315	rVB2	2353	6027	1.07%	0.129%
4	5.501	702	713	727	rBV	70117	199158	35.23%	4.259%
5	5.659	727	739	756	rVB	130085	364252	64.44%	7.790%
6	6.061	793	805	823	rBV2	77910	208546	36.89%	4.460%
7	6.860	924	936	954	rBV2	182925	425633	75.30%	9.102%
8	8.713	1233	1240	1255	rBV	356814	565251	100.00%	12.088%
9	10.116	1464	1470	1482	rBV	355724	489340	86.57%	10.465%
10	10.359	1505	1510	1516	rBV	7598	10520	1.86%	0.225%
11	10.707	1561	1567	1574	rBV3	3933	7356	1.30%	0.157%
12	11.140	1632	1638	1652	rBV	313017	399585	70.69%	8.545%
13	11.432	1682	1686	1689	rBV	14389	22662	4.01%	0.485%
14	11.506	1695	1698	1707	rVB	4699	6469	1.14%	0.138%
15	11.670	1720	1725	1730	rBV2	5148	6887	1.22%	0.147%
16	11.810	1743	1748	1753	rBV	48789	60165	10.64%	1.287%
17	12.079	1787	1792	1799	rBV	439681	531872	94.09%	11.375%
18	12.146	1799	1803	1807	rVB	15072	18825	3.33%	0.403%
19	12.298	1822	1828	1830	rBV	19028	24696	4.37%	0.528%
20	12.322	1830	1832	1836	rVV	20423	23412	4.14%	0.501%
21	12.371	1836	1840	1848	rVB3	23994	42239	7.47%	0.903%
22	12.469	1851	1856	1860	rBV	11932	14655	2.59%	0.313%
23	12.518	1860	1864	1869	rVB	7227	8861	1.57%	0.189%
24	12.585	1871	1875	1877	rBV	11457	14220	2.52%	0.304%
25	12.609	1877	1879	1884	rVB	15000	16876	2.99%	0.361%
26	12.670	1884	1889	1893	rBV	36259	43472	7.69%	0.930%
27	12.774	1900	1906	1913	rBV5	2791	6617	1.17%	0.142%
28	12.908	1923	1928	1932	rVB	6598	8325	1.47%	0.178%
29	12.975	1932	1939	1943	rBV	15434	20004	3.54%	0.428%
30	13.017	1943	1946	1952	rVB	19223	23858	4.22%	0.510%
31	13.225	1974	1980	1984	rBV2	15826	20969	3.71%	0.448%
32	13.347	1996	2000	2005	rVB2	18326	25148	4.45%	0.538%
33	13.402	2005	2009	2012	rBV2	11146	13542	2.40%	0.290%
34	13.481	2018	2022	2028	rVB	13879	19060	3.37%	0.408%

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Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	13.560	2028	2035	2044	rBV	5756	9102	1.61%	0.195%
36	13.633	2044	2047	2054	rVB4	3454	6154	1.09%	0.132%
37	13.834	2074	2080	2088	rBV	469442	561330	99.31%	12.004%
38	14.383	2163	2170	2176	rBV4	3333	7378	1.31%	0.158%
39	14.694	2214	2221	2231	rBV	171619	221278	39.15%	4.732%
40	14.840	2239	2245	2253	rVB	73524	92269	16.32%	1.973%
41	15.273	2308	2316	2322	rBV	15959	23016	4.07%	0.492%
42	15.444	2338	2344	2347	rBV	6650	9435	1.67%	0.202%
43	15.547	2356	2361	2366	rBV	9152	15064	2.67%	0.322%
44	15.688	2378	2384	2388	rBV	9257	17122	3.03%	0.366%
45	15.730	2388	2391	2397	rVB2	6394	9519	1.68%	0.204%
46	15.913	2414	2421	2429	rBV5	2950	7360	1.30%	0.157%

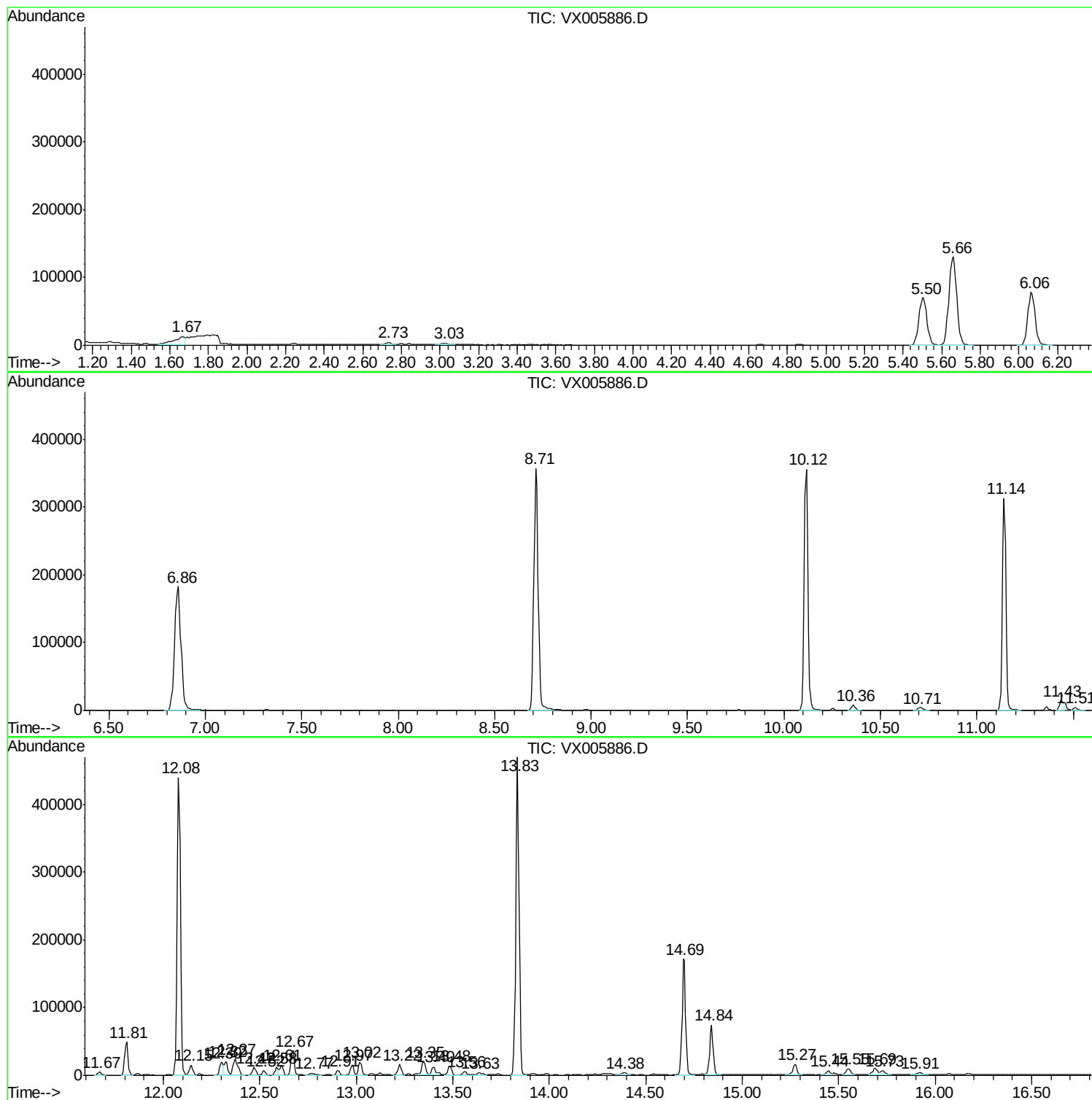
Sum of corrected areas: 4676003

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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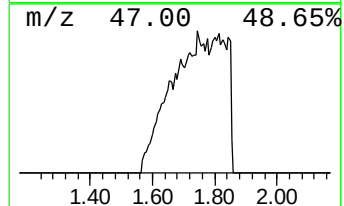
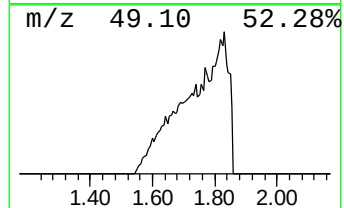
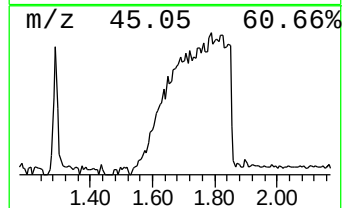
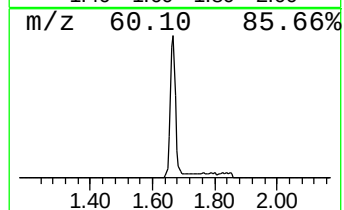
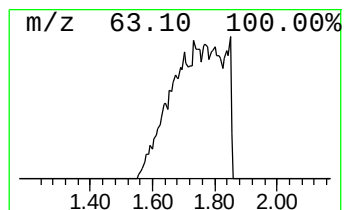
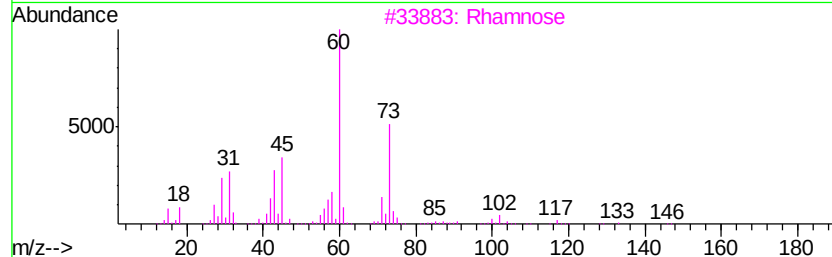
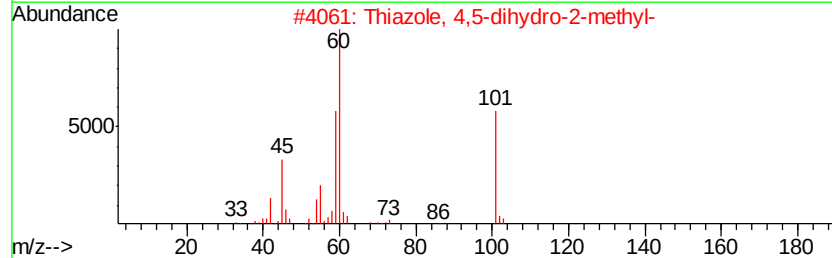
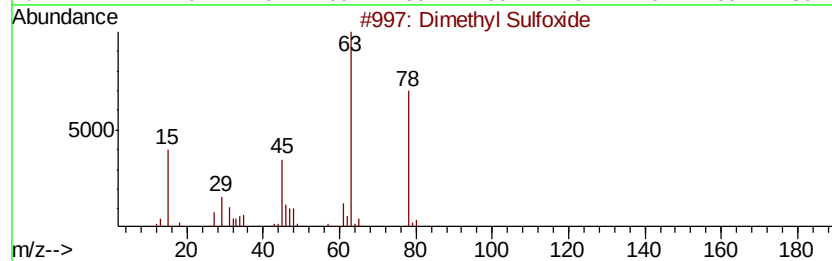
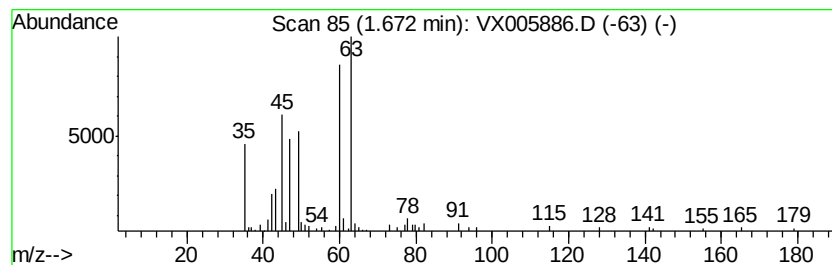
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown1.67 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.67	5.76 ug/l	41936	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl Sulfoxide	78	C2H6OS	000067-68-5	12
2		Thiazole, 4,5-dihydro-2-methyl-	101	C4H7NS	002346-00-1	10
3		Rhamnose	164	C6H12O5	003615-41-6	9
4		Heptanoic acid	130	C7H14O2	000111-14-8	9
5		Acetamide, N-(.beta.-mercaptoeth...	119	C4H9NOS	001190-73-4	9



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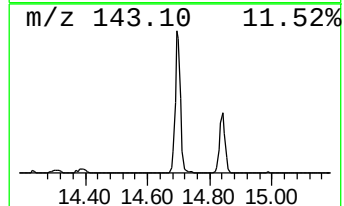
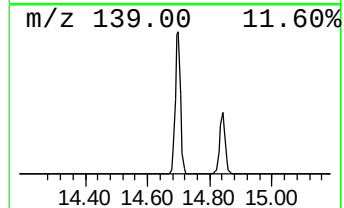
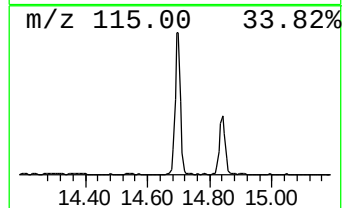
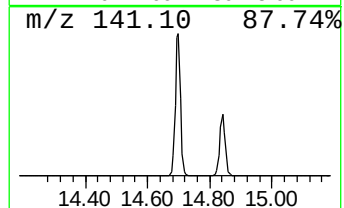
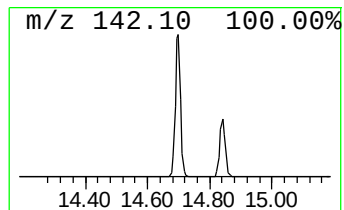
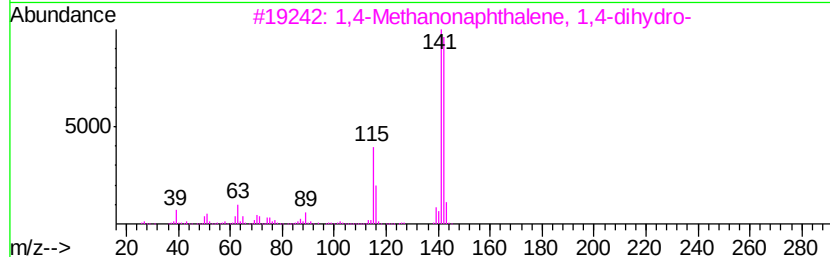
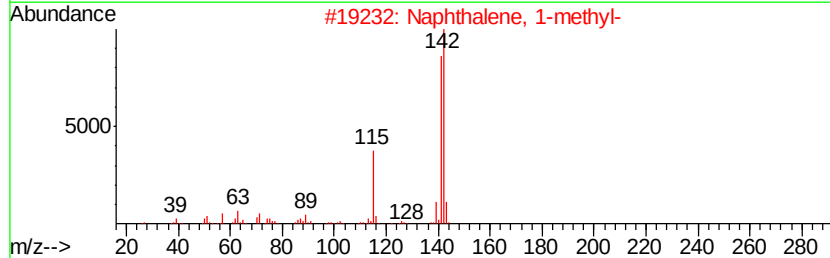
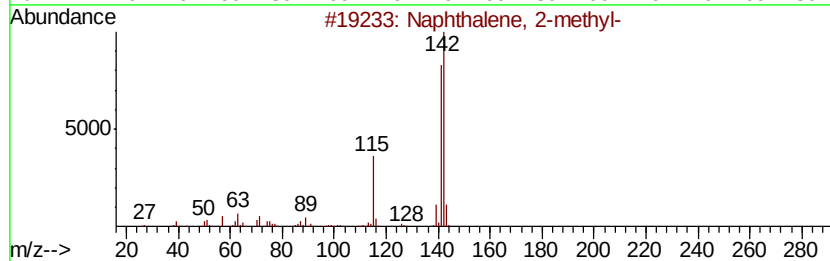
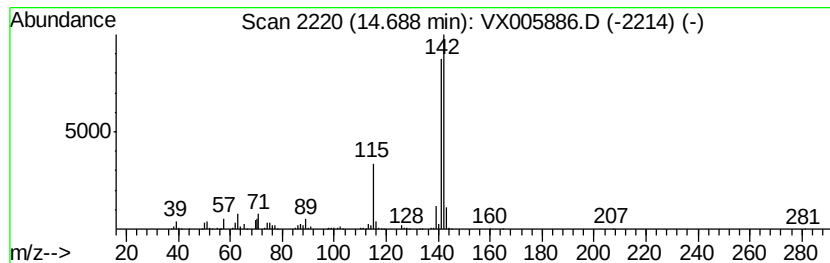
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Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	20.80 ug/l	221278	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		3-Indolecarbonitrile	142 C9H6N2	005457-28-3 46



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.67	1.67	5.8	ug/l	41936	1	5.66	364252	50.0
Naphthalene, 1-me...	14.69	20.8	ug/l	221278	4	12.08	531872	50.0