

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110918\
 Data File : VX005888.D
 Acq On : 10 Nov 2018 03:58
 Operator : JC/MD
 Sample : J5889-14
 Misc : 13.32g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BOTTOM-E

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	65	84	87	rBV2	13047	47335	8.25%	1.083%
2	3.019	300	306	316	rBV	2977	6488	1.13%	0.148%
3	5.501	702	713	727	rBV	72820	205186	35.76%	4.696%
4	5.659	728	739	757	rVB2	134283	366457	63.87%	8.387%
5	6.068	795	806	825	rBV2	78384	212988	37.12%	4.875%
6	6.860	926	936	958	rBV	190747	445565	77.66%	10.198%
7	8.713	1233	1240	1258	rBV	348497	573757	100.00%	13.131%
8	10.116	1464	1470	1481	rBV	352330	492189	85.78%	11.265%
9	10.359	1505	1510	1515	rVB	6198	8701	1.52%	0.199%
10	10.701	1562	1566	1573	rBV3	3085	6384	1.11%	0.146%
11	11.140	1632	1638	1653	rBV	307086	394703	68.79%	9.034%
12	11.439	1682	1687	1689	rBV	12182	18482	3.22%	0.423%
13	11.506	1695	1698	1709	rVB3	3669	6475	1.13%	0.148%
14	11.810	1743	1748	1754	rVB	38396	46656	8.13%	1.068%
15	12.079	1787	1792	1799	rBV	415090	518004	90.28%	11.855%
16	12.146	1799	1803	1807	rVB	11375	14862	2.59%	0.340%
17	12.304	1824	1829	1830	rBV	13836	17173	2.99%	0.393%
18	12.323	1830	1832	1836	rVB	15879	16986	2.96%	0.389%
19	12.371	1836	1840	1848	rVB3	19417	33434	5.83%	0.765%
20	12.475	1851	1857	1861	rBV2	10548	14823	2.58%	0.339%
21	12.518	1861	1864	1870	rVB	5434	7625	1.33%	0.175%
22	12.591	1870	1876	1877	rBV	9026	11639	2.03%	0.266%
23	12.609	1877	1879	1884	rVB	11406	13484	2.35%	0.309%
24	12.670	1884	1889	1893	rBV	27728	32724	5.70%	0.749%
25	12.908	1924	1928	1932	rVB	5091	6468	1.13%	0.148%
26	12.981	1932	1940	1943	rBV	11761	17084	2.98%	0.391%
27	13.018	1943	1946	1954	rVB	14264	19462	3.39%	0.445%
28	13.225	1975	1980	1984	rBV2	12522	16627	2.90%	0.381%
29	13.322	1991	1996	1997	rBV3	9164	11662	2.03%	0.267%
30	13.347	1997	2000	2005	rVB2	15648	22079	3.85%	0.505%
31	13.402	2005	2009	2012	rBV2	7933	10444	1.82%	0.239%
32	13.481	2018	2022	2030	rVB	11345	17265	3.01%	0.395%
33	13.560	2031	2035	2041	rVV2	4351	5938	1.03%	0.136%
34	13.834	2075	2080	2087	rVB	345146	414600	72.26%	9.489%

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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

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35	13.902	2087	2091	2097	rVB5	4172	6528	1.14%	0.149%
36	14.091	2115	2122	2125	rBV	9113	11448	2.00%	0.262%
37	14.694	2214	2221	2231	rBV	118228	159055	27.72%	3.640%
38	14.810	2236	2240	2241	rVV	7772	7384	1.29%	0.169%
39	14.840	2241	2245	2251	rVB	51726	65610	11.44%	1.502%
40	15.273	2309	2316	2320	rBV2	11545	18251	3.18%	0.418%
41	15.444	2340	2344	2347	rBV2	4368	6279	1.09%	0.144%
42	15.487	2348	2351	2356	rVB4	4792	7323	1.28%	0.168%
43	15.554	2356	2362	2367	rBV2	8473	14621	2.55%	0.335%
44	15.688	2379	2384	2388	rBV2	7089	10771	1.88%	0.247%
45	15.822	2400	2406	2413	rVB9	4025	8304	1.45%	0.190%

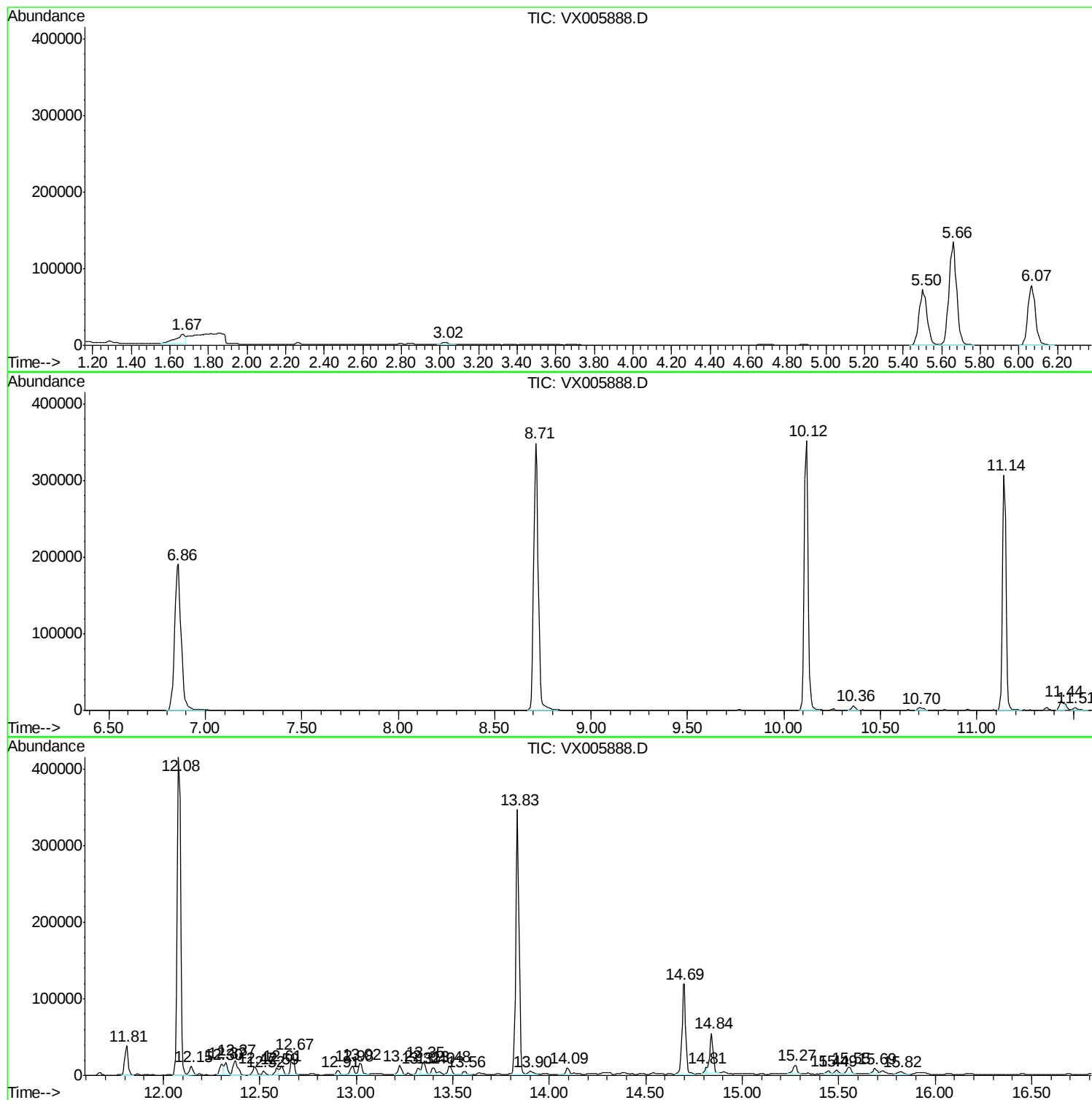
Sum of corrected areas: 4369323

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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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MSVOA_X
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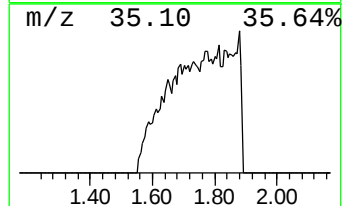
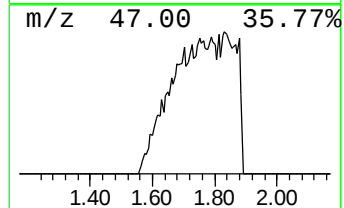
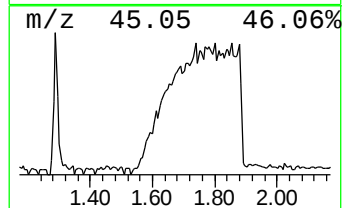
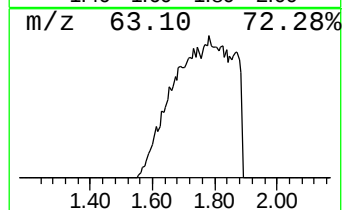
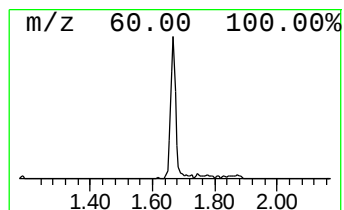
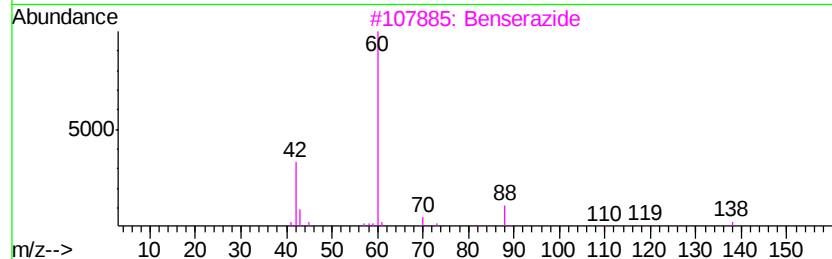
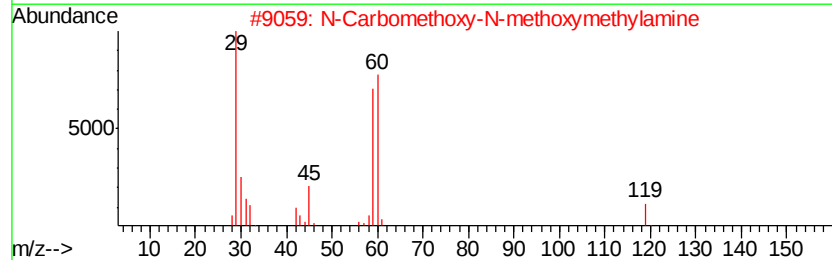
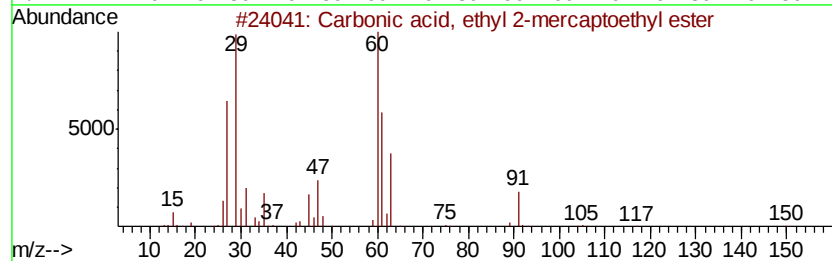
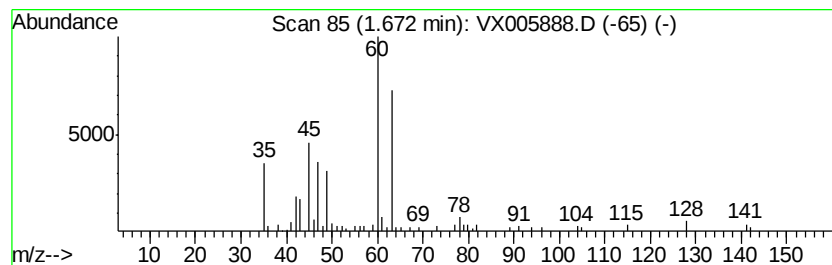
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown1.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.67	6.46 ug/l	47335	Pentafluorobenzene	5.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, ethyl 2-mercaptoe...	150	C5H10O3S	005628-96-6	28
2			N-Carbomethoxy-N-methoxymethylamine	119	C4H9NO3	153654-07-0	9
3			Benserazide	257	C10H15N3O5	000322-35-0	9
4			1,2-Ethanediol, diformate	118	C4H6O4	000629-15-2	9
5			1,2-Ethanedithiol	94	C2H6S2	000540-63-6	9



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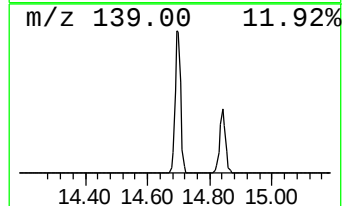
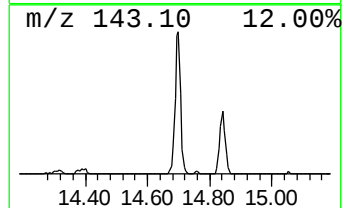
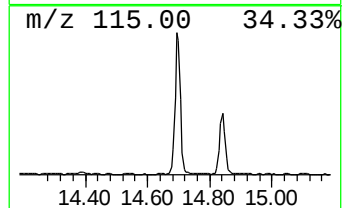
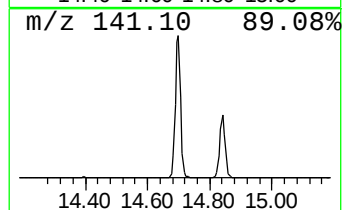
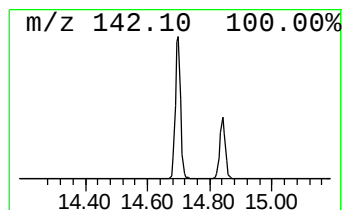
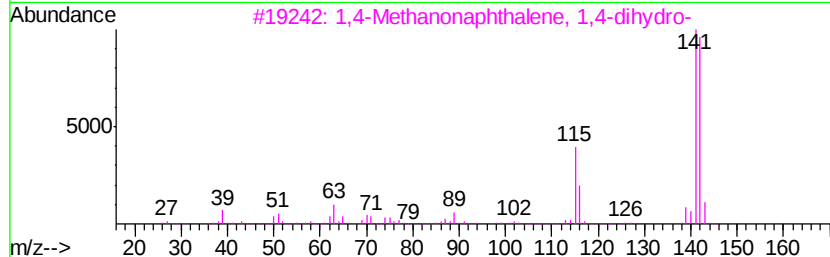
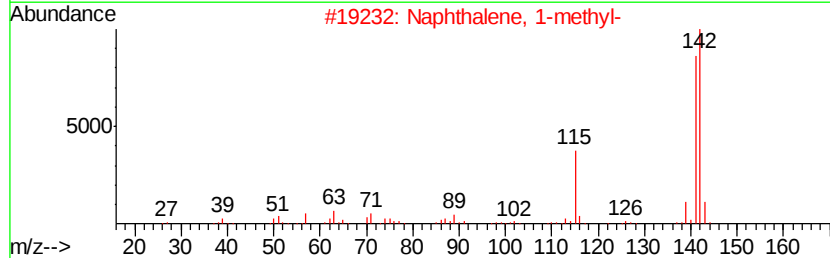
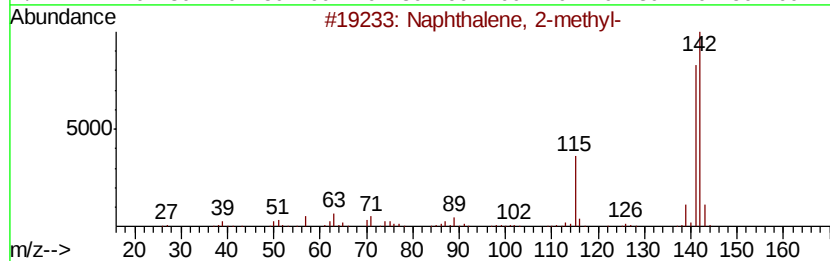
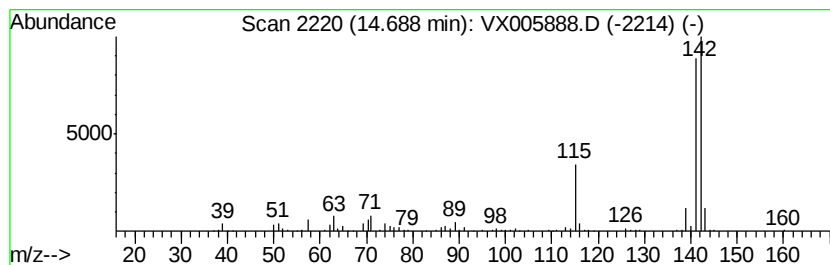
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Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	15.35 ug/l	159055	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 90
4		Benzocycloheptatriene	142 C11H10	000264-09-5 90
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.67	1.67	6.5	ug/l	47335	1	5.66	366457	50.0
Naphthalene, 1-me...	14.69	15.4	ug/l	159055	4	12.08	518004	50.0