

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
 Data File : VX017547.D
 Acq On : 24 Jul 2020 16:53
 Operator : JC/SP
 Sample : L3413-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 20758

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\82X072320W.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	2.416	212	218	231	rBV	4633106	7821848	72.90%	6.022%
2	2.581	237	245	261	rBV	2402861	4217472	39.31%	3.247%
3	3.166	332	341	351	rBV	148171	353433	3.29%	0.272%
4	4.398	530	543	554	rBV2	75441	244795	2.28%	0.188%
5	4.635	571	582	595	rBV	202148	585460	5.46%	0.451%
6	4.794	599	608	620	rVB2	364719	1021735	9.52%	0.787%
7	5.086	647	656	666	rBV2	119776	384832	3.59%	0.296%
8	5.464	706	718	729	rBV	292026	851911	7.94%	0.656%
9	5.629	734	745	757	rBV	447613	1256940	11.71%	0.968%
10	6.037	802	812	817	rBV	310116	778977	7.26%	0.600%
11	6.117	817	825	833	rVV	876493	2325392	21.67%	1.790%
12	6.178	833	835	847	rVB2	44866	119063	1.11%	0.092%
13	6.830	933	942	959	rVB	692511	1662780	15.50%	1.280%
14	7.519	1048	1055	1065	rVB2	95765	197293	1.84%	0.152%
15	8.695	1241	1248	1254	rBV	1728064	2573583	23.98%	1.981%
16	8.769	1254	1260	1265	rVV	6998903	10730070	100.00%	8.261%
17	8.811	1265	1267	1277	rVB	162814	252365	2.35%	0.194%
18	10.098	1467	1478	1489	rBV	1622288	2229708	20.78%	1.717%
19	10.238	1495	1501	1507	rBV	1102205	1472677	13.72%	1.134%
20	10.342	1512	1518	1525	rVB	5292584	6745529	62.87%	5.193%
21	10.683	1568	1574	1585	rBV	5326655	6693730	62.38%	5.153%
22	11.000	1620	1626	1630	rBV	95007	130163	1.21%	0.100%
23	11.122	1640	1646	1656	rBV	1629011	2031388	18.93%	1.564%
24	11.341	1678	1682	1689	rVB	216415	273830	2.55%	0.211%
25	11.421	1689	1695	1702	rVV	1657211	2693295	25.10%	2.074%
26	11.488	1702	1706	1719	rVB	910420	1321431	12.32%	1.017%
27	11.652	1728	1733	1742	rBV	1287751	1539149	14.34%	1.185%
28	11.793	1751	1756	1762	rVB	5663050	6667813	62.14%	5.133%
29	11.933	1775	1779	1784	rVB	78284	108616	1.01%	0.084%
30	12.012	1788	1792	1795	rVB	101352	125424	1.17%	0.097%
31	12.061	1795	1800	1806	rBV	2361453	2902816	27.05%	2.235%
32	12.128	1806	1811	1816	rBV	2753308	3505181	32.67%	2.699%
33	12.280	1828	1836	1844	rBV2	2900963	4473399	41.69%	3.444%
34	12.359	1844	1849	1854	rVB3	613407	947818	8.83%	0.730%

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35	12.457	1854	1865	1869	rBV3	347753	563072	5.25%	0.433%
36	12.500	1869	1872	1878	rVB	274514	359893	3.35%	0.277%
37	12.573	1878	1884	1886	rBV	514977	732204	6.82%	0.564%
38	12.591	1886	1887	1892	rVB	477978	449826	4.19%	0.346%
39	12.652	1892	1897	1900	rBV	726974	828233	7.72%	0.638%
40	12.683	1900	1902	1906	rVB	196059	214862	2.00%	0.165%
41	12.737	1906	1911	1919	rBV2	728524	1179050	10.99%	0.908%
42	12.835	1923	1927	1932	rBV	147969	190372	1.77%	0.147%
43	12.890	1932	1936	1940	rVB	343899	421930	3.93%	0.325%
44	12.963	1940	1948	1951	rBV	615875	878587	8.19%	0.676%
45	13.000	1951	1954	1960	rVB	1103430	1328232	12.38%	1.023%
46	13.085	1960	1968	1970	rBV4	154705	331092	3.09%	0.255%
47	13.207	1980	1988	1992	rBV	1347955	1787245	16.66%	1.376%
48	13.250	1992	1995	1998	rVB	173799	183981	1.71%	0.142%
49	13.329	2004	2008	2013	rVB2	3008973	4191667	39.06%	3.227%
50	13.378	2013	2016	2019	rVB3	89841	112906	1.05%	0.087%
51	13.463	2025	2030	2038	rVB	2012800	2533759	23.61%	1.951%
52	13.542	2038	2043	2046	rBV2	257777	372760	3.47%	0.287%
53	13.585	2046	2050	2053	rBV	363191	457042	4.26%	0.352%
54	13.621	2053	2056	2061	rVV3	493530	833668	7.77%	0.642%
55	13.676	2062	2065	2067	rVV	192604	283507	2.64%	0.218%
56	13.707	2067	2070	2074	rVB2	528656	748454	6.98%	0.576%
57	13.817	2083	2088	2094	rVB	4441920	5545129	51.68%	4.269%
58	13.896	2094	2101	2106	rVB2	831131	1224545	11.41%	0.943%
59	13.963	2106	2112	2117	rBV2	847196	1308273	12.19%	1.007%
60	14.012	2117	2120	2124	rVV	270090	295585	2.75%	0.228%
61	14.079	2127	2131	2133	rVV3	323648	368153	3.43%	0.283%
62	14.115	2133	2137	2141	rVB	721856	798082	7.44%	0.614%
63	14.268	2155	2162	2168	rVB	2445652	3694260	34.43%	2.844%
64	14.359	2174	2177	2182	rBV	242786	325015	3.03%	0.250%
65	14.414	2182	2186	2190	rVV	775622	943439	8.79%	0.726%
66	14.457	2190	2193	2198	rVB2	166399	265518	2.47%	0.204%
67	14.518	2198	2203	2213	rBV2	1634083	2414056	22.50%	1.859%
68	14.597	2213	2216	2222	rVB6	55575	110275	1.03%	0.085%
69	14.676	2222	2229	2233	rBV2	2128304	3684702	34.34%	2.837%
70	14.713	2233	2235	2240	rVB2	506862	579013	5.40%	0.446%
71	14.768	2240	2244	2246	rBV3	199104	289151	2.69%	0.223%

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72	14.792	2246	2248	2250	rVV	529807	608699	5.67%	0.469%
73	14.822	2250	2253	2256	rVV	1200723	1520603	14.17%	1.171%
74	14.859	2256	2259	2261	rVV2	282151	395698	3.69%	0.305%
75	14.890	2261	2264	2269	rVV4	290319	390705	3.64%	0.301%
76	14.944	2269	2273	2280	rVB2	312383	527768	4.92%	0.406%
77	15.097	2291	2298	2302	rBV4	201107	400965	3.74%	0.309%
78	15.231	2315	2320	2322	rBV	713036	1104369	10.29%	0.850%
79	15.304	2327	2332	2333	rBV4	85295	118783	1.11%	0.091%
80	15.420	2347	2351	2355	rBV2	280275	440410	4.10%	0.339%
81	15.536	2364	2370	2374	rBV2	816700	1352189	12.60%	1.041%
82	15.664	2386	2391	2395	rVV	554208	1012983	9.44%	0.780%
83	15.700	2395	2397	2405	rVB	325345	479721	4.47%	0.369%
84	15.895	2425	2429	2433	rBV3	135820	221336	2.06%	0.170%
85	16.042	2450	2453	2457	rBV4	91676	141394	1.32%	0.109%
86	16.078	2457	2459	2464	rVB4	97938	144421	1.35%	0.111%
87	16.145	2465	2470	2478	rBV3	196296	415529	3.87%	0.320%
88	16.432	2512	2517	2524	rVB	316770	547154	5.10%	0.421%

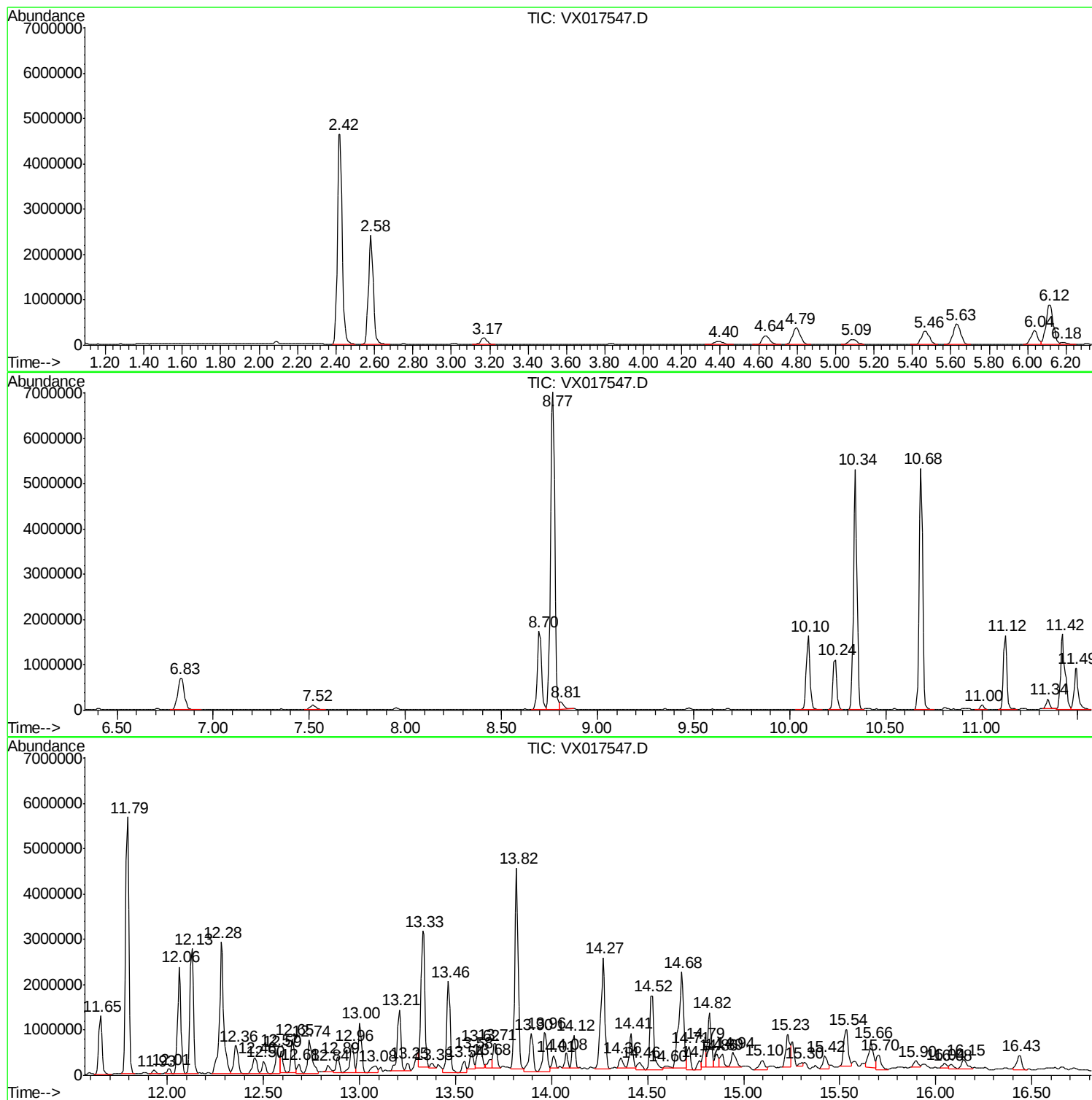
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260

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



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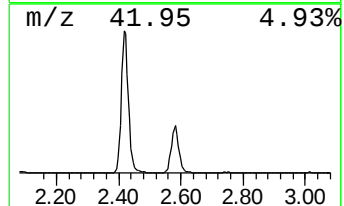
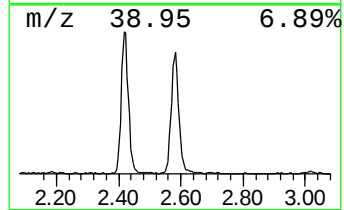
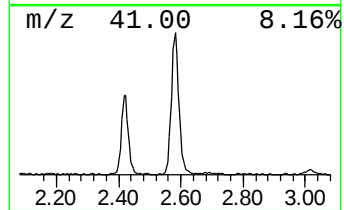
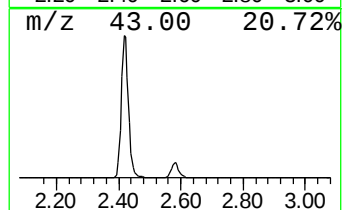
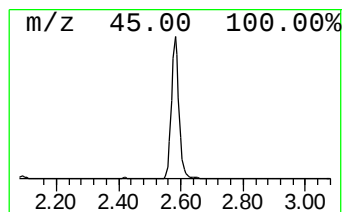
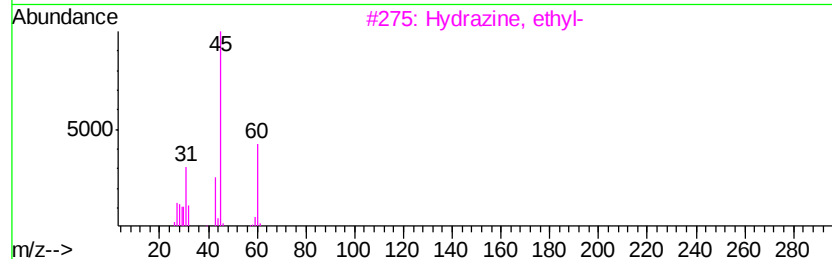
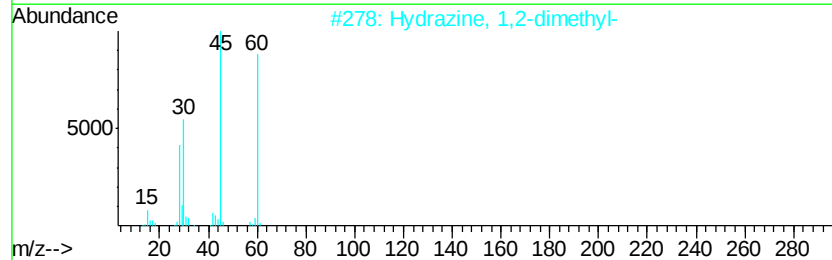
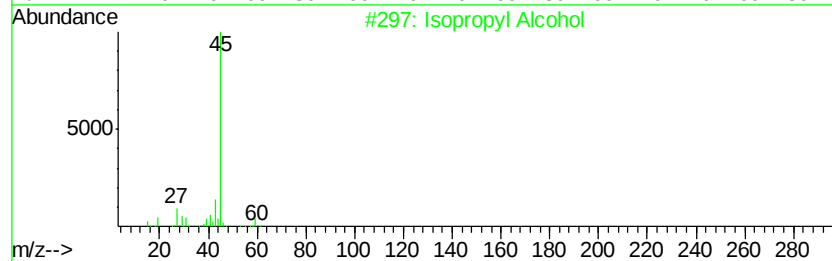
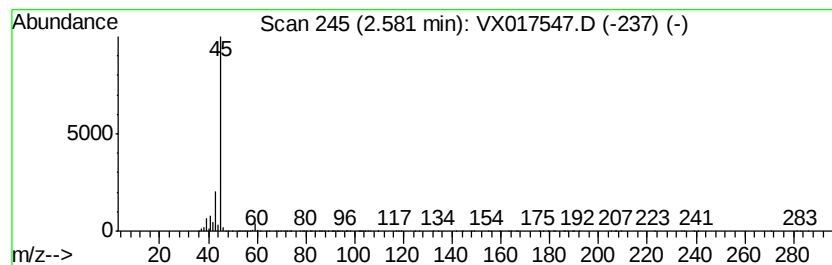
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260

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

Peak Number 1 Isopropyl Alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.58	167.77 ug/l	4217470	Pentafluorobenzene	5.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isopropyl Alcohol	60	C3H8O	000067-63-0	86
2		Hydrazine, 1,2-dimethyl-	60	C2H8N2	000540-73-8	9
3		Hydrazine, ethyl-	60	C2H8N2	000624-80-6	9
4		Formamide	45	CH3NO	000075-12-7	5
5		Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	4



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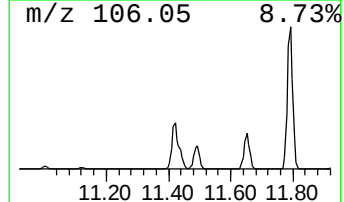
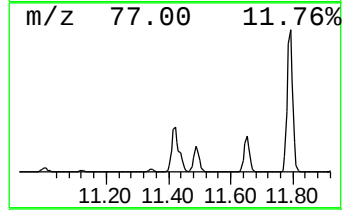
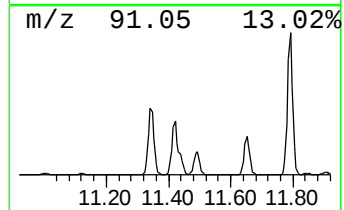
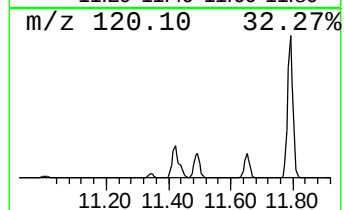
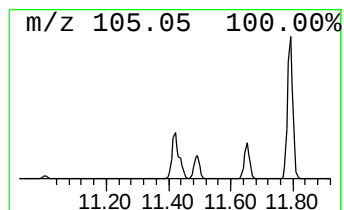
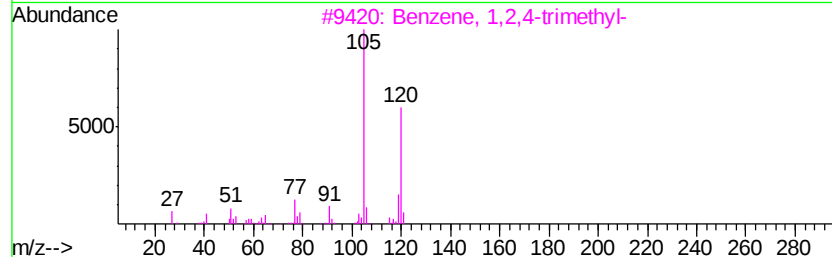
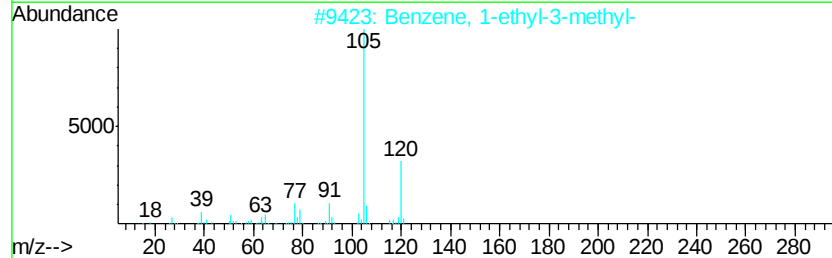
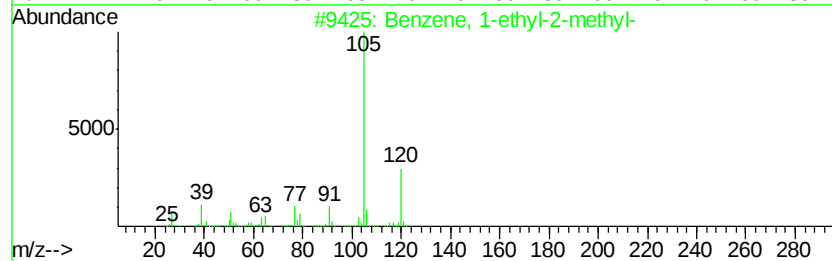
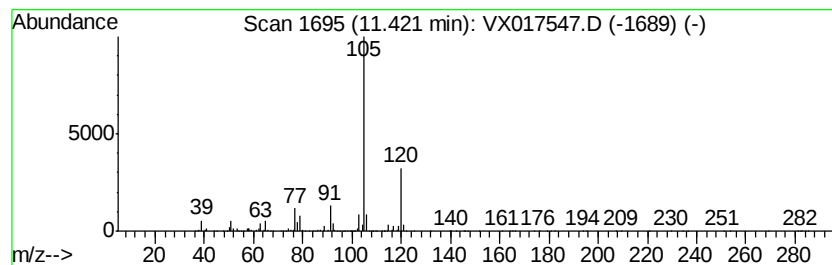
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	46.39 ug/l	2693300	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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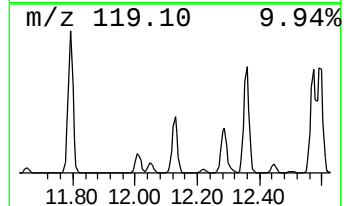
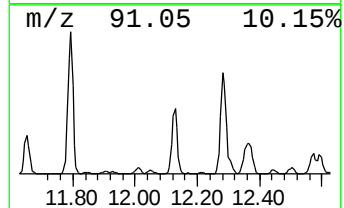
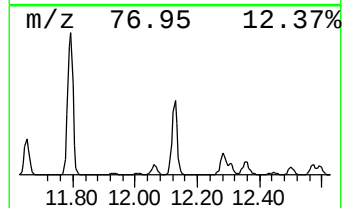
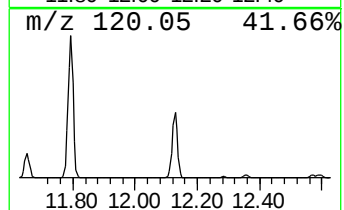
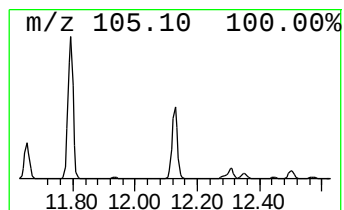
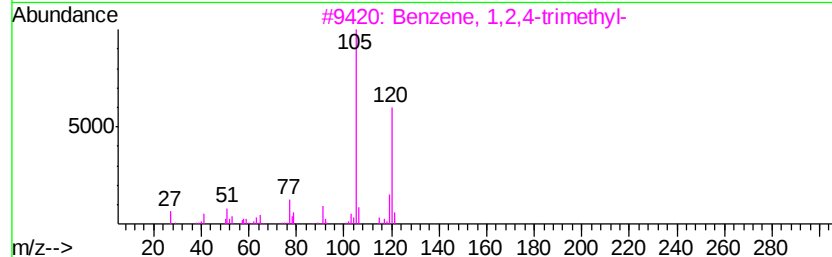
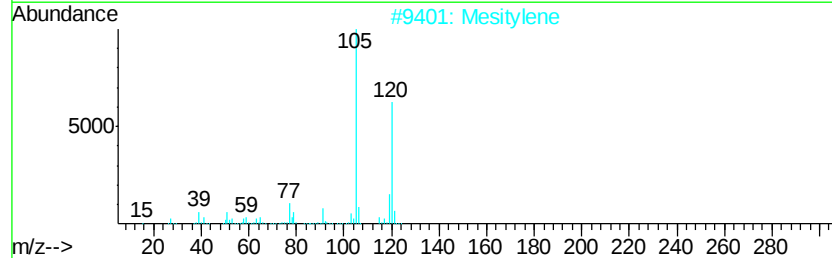
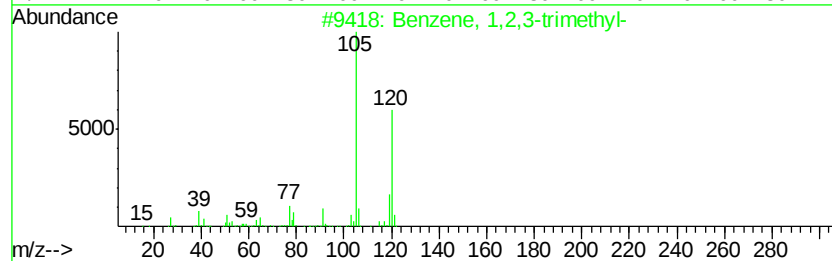
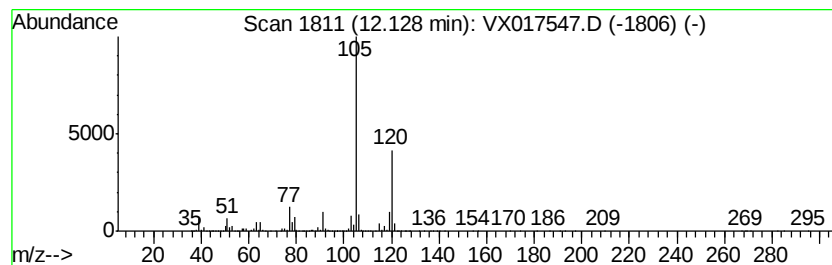
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.13	60.38 ug/l	3505180	1,4-Dichlorobenzene-d4	12.06	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8	95
2		Mesitylene	120 C9H12	000108-67-8	91
3		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8	90



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20758

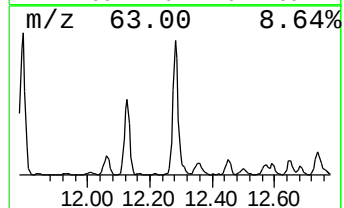
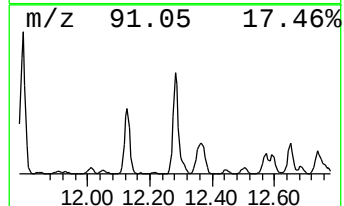
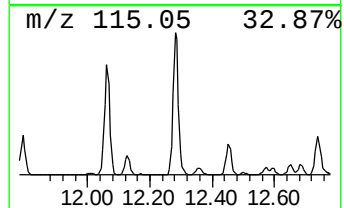
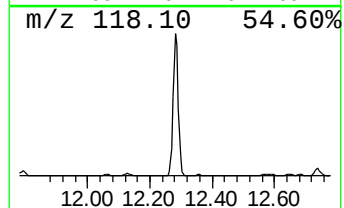
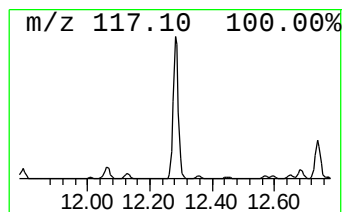
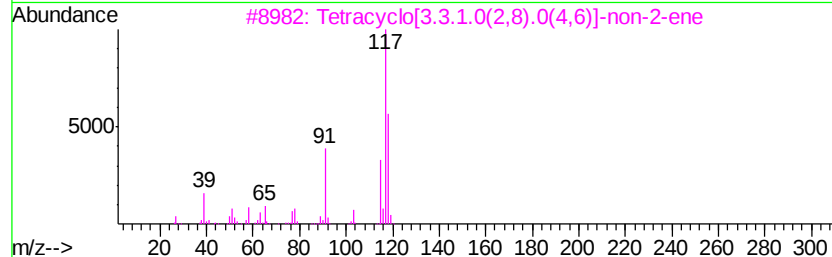
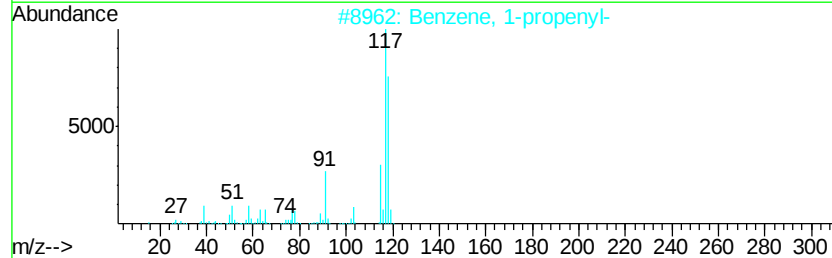
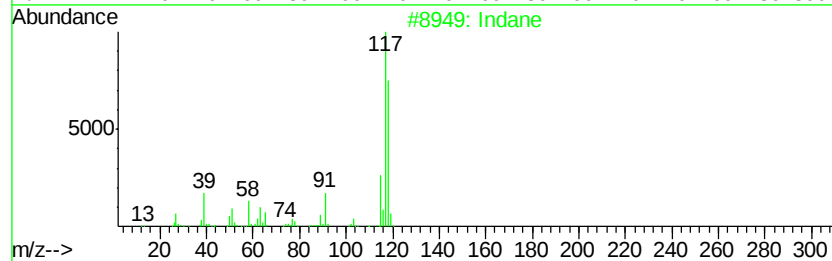
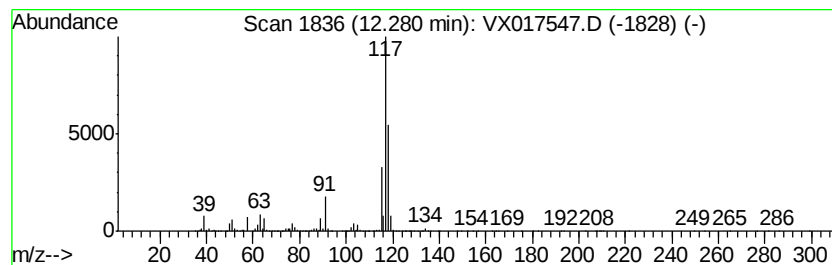
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	77.05 ug/l	4473400	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	91
2		Benzene, 1-propenyl-	118	C9H10	000637-50-3	81
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
20758

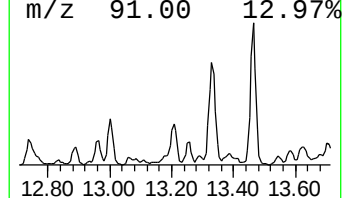
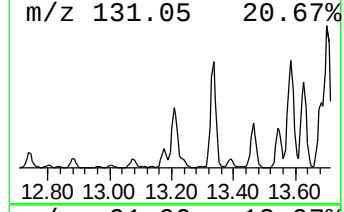
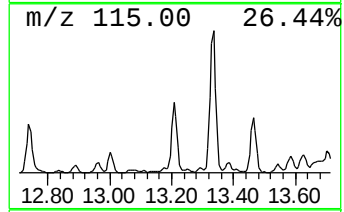
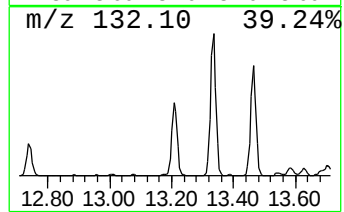
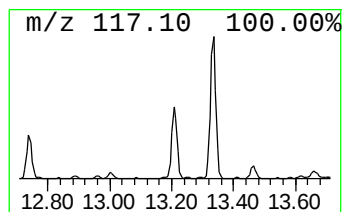
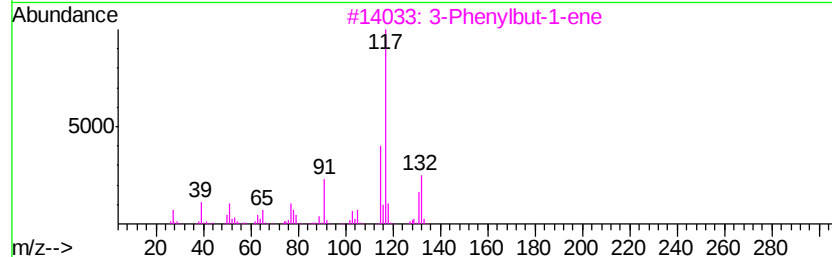
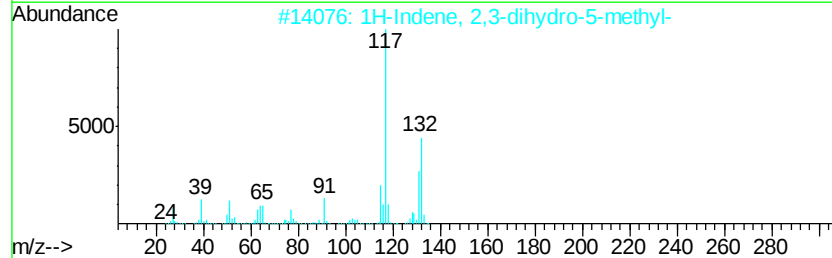
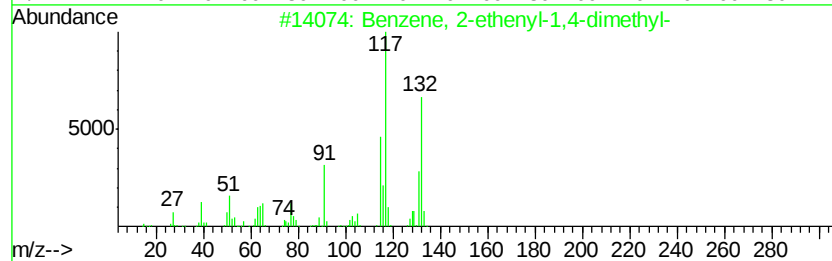
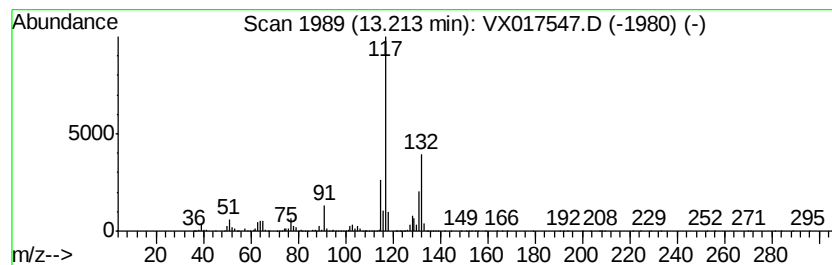
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	30.78 ug/l	1787250	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
20758

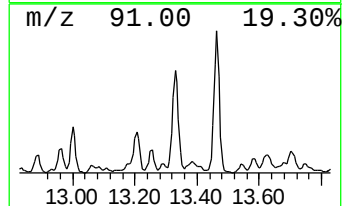
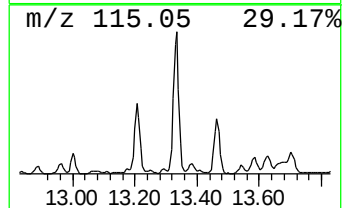
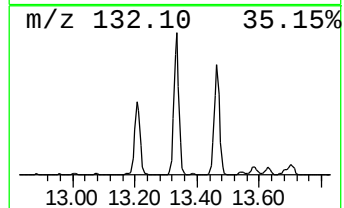
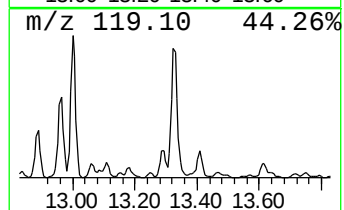
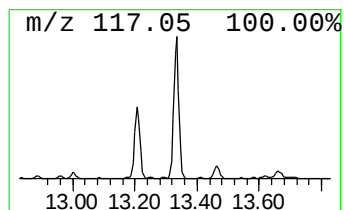
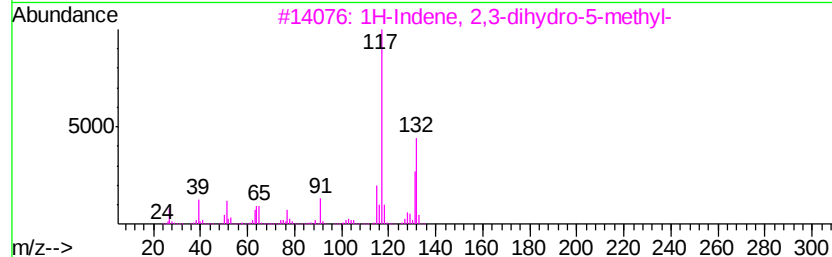
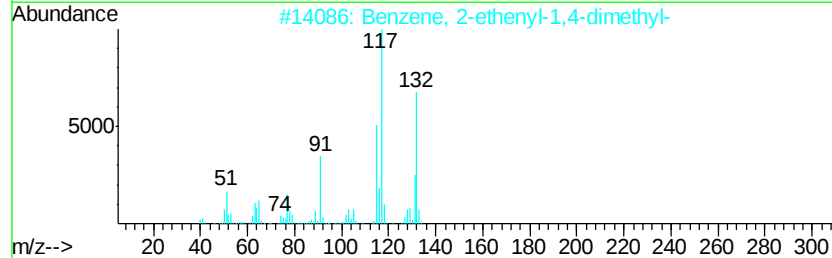
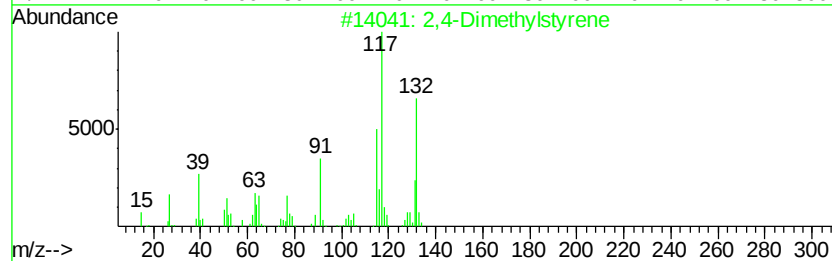
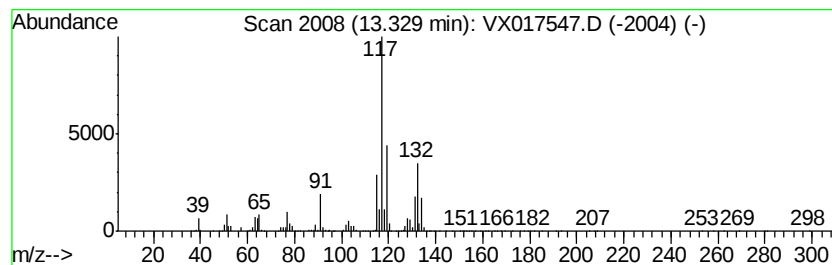
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Quant Title : SW846 8260

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

Peak Number 6 2,4-Dimethylstyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	72.20 ug/l	4191670	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	76
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
20758

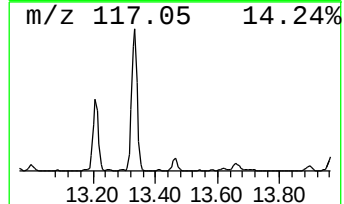
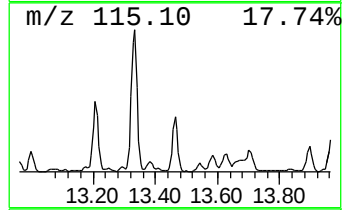
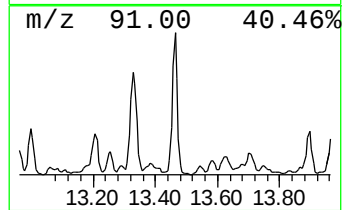
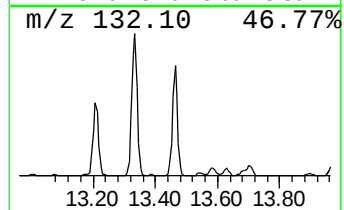
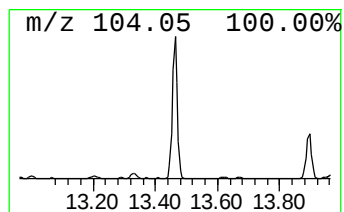
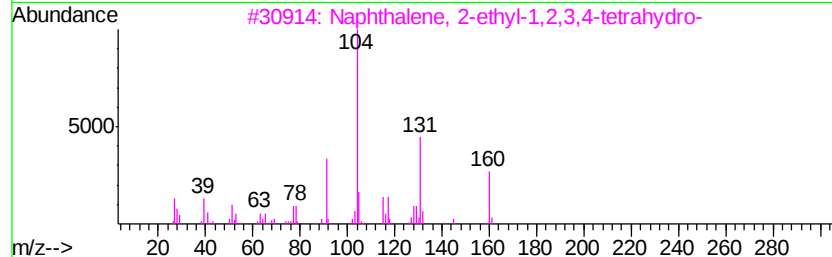
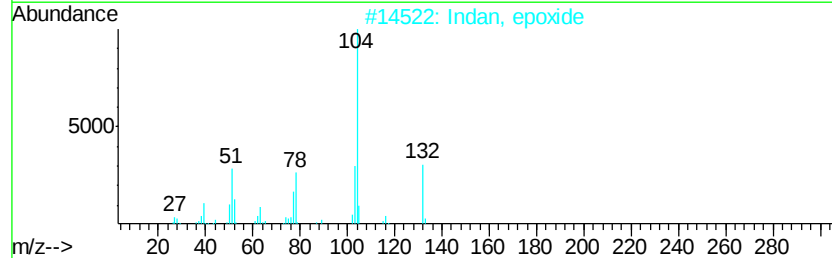
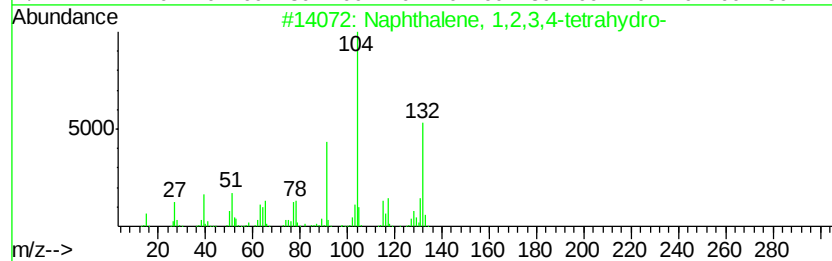
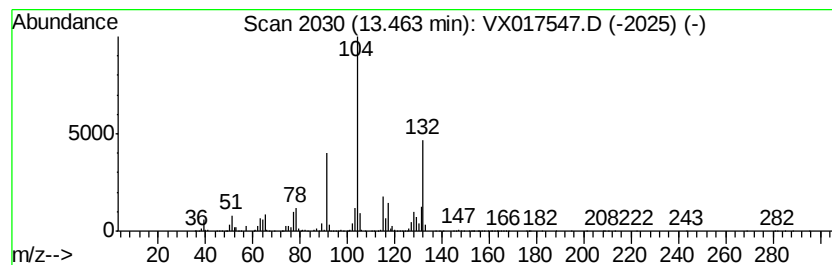
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Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	43.64 ug/l	2533760	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	91
2		Indan, epoxide	132	C9H8O	000768-22-9	58
3		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	53
4		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5		Benzene, cyclobutyl-	132	C10H12	004392-30-7	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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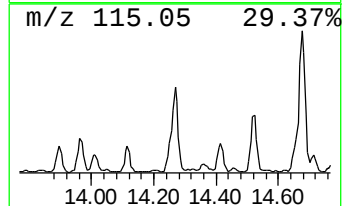
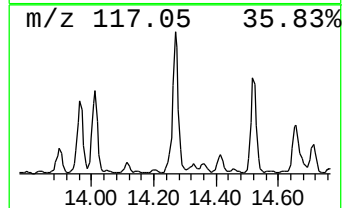
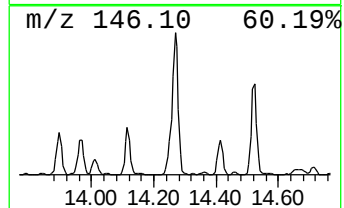
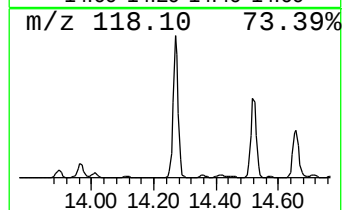
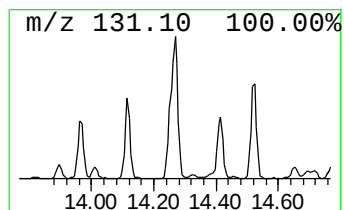
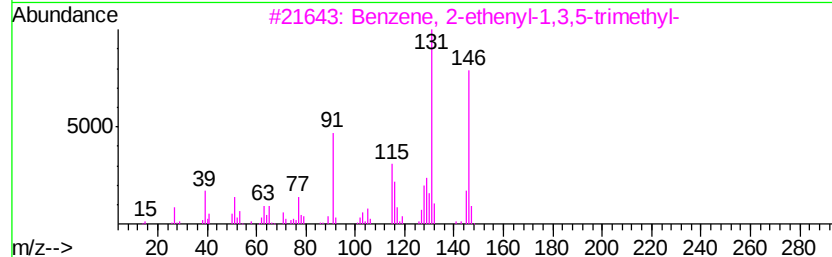
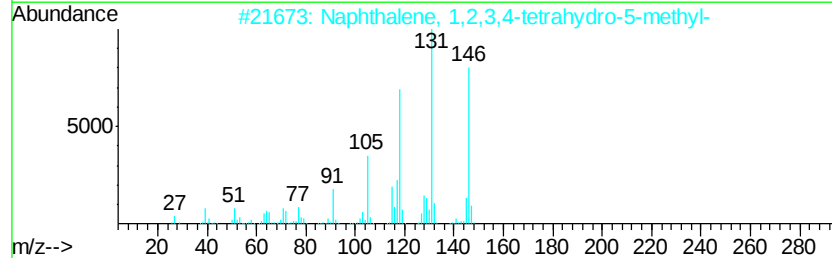
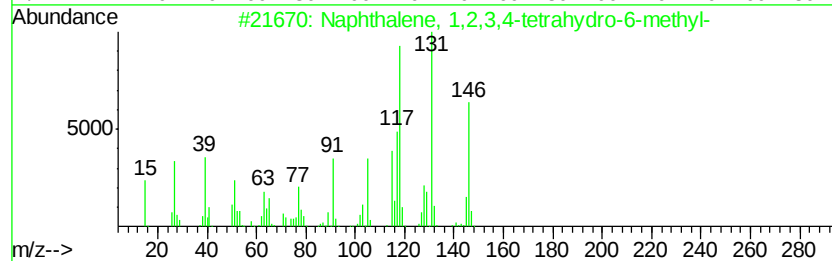
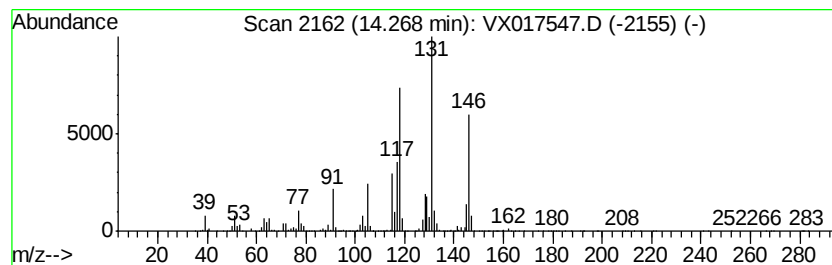
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	63.63 ug/l	3694260	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
4		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	64
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
20758

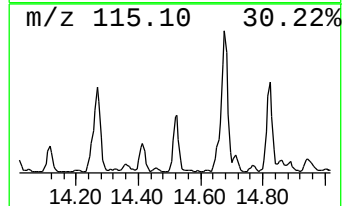
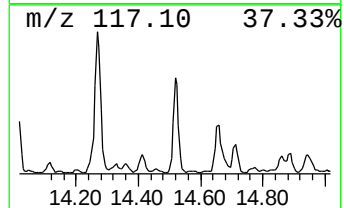
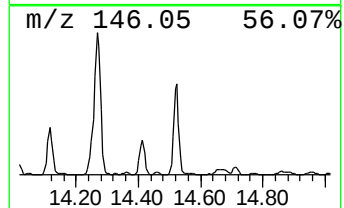
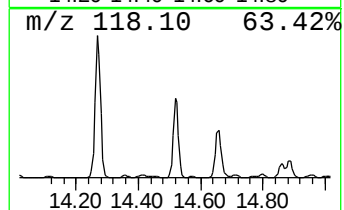
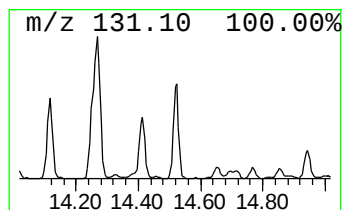
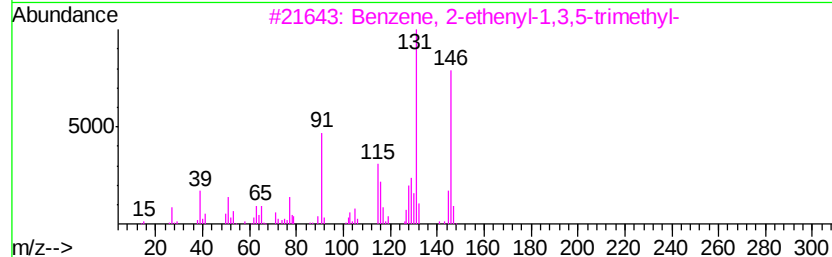
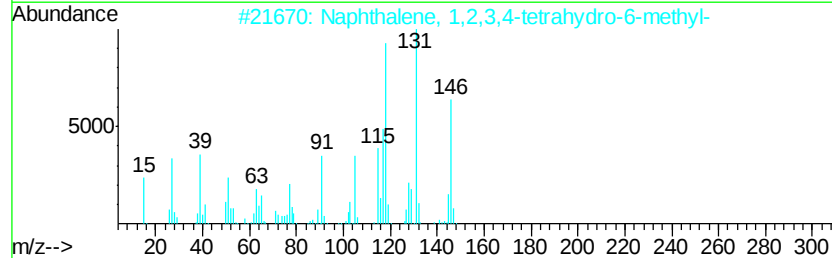
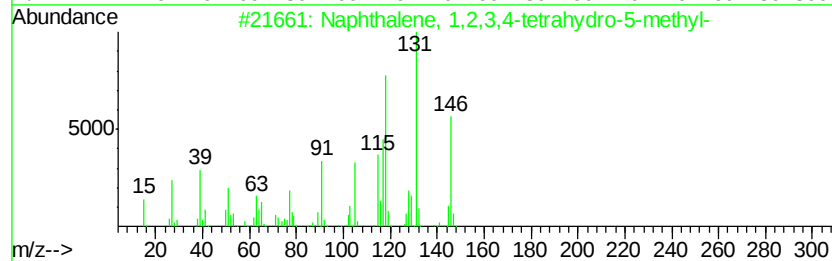
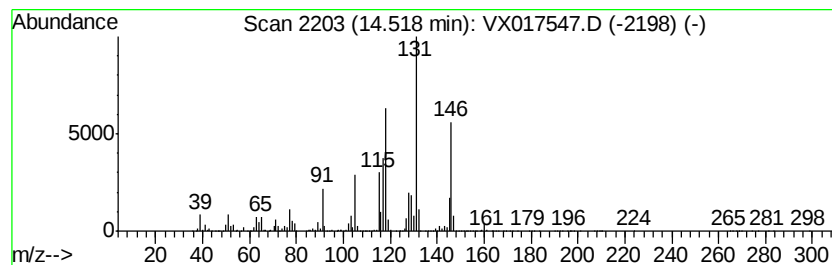
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	41.58 ug/l	2414060	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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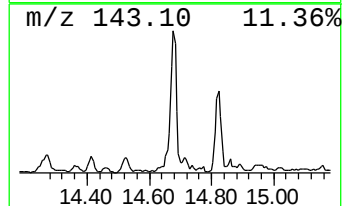
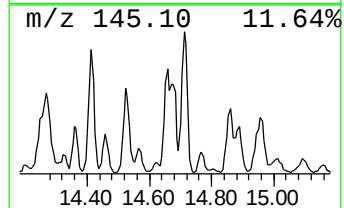
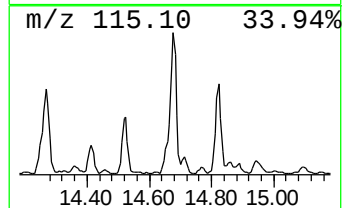
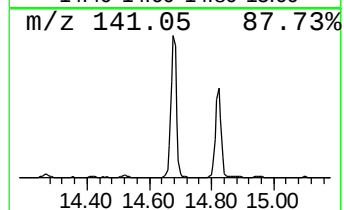
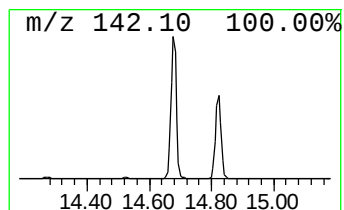
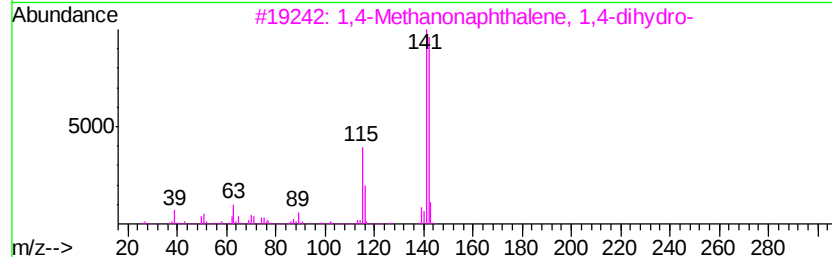
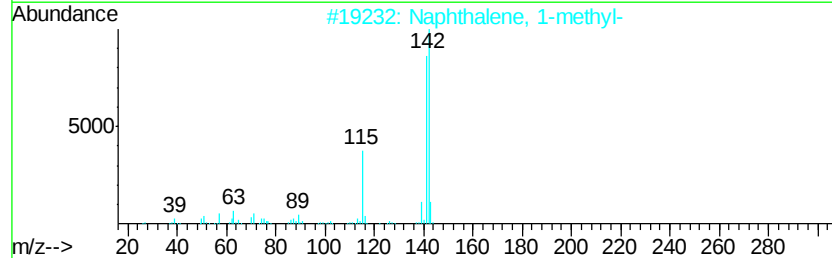
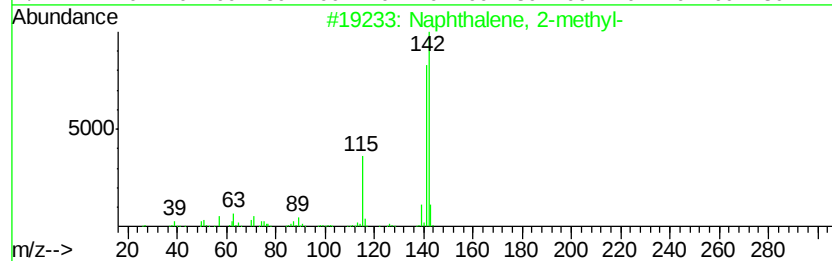
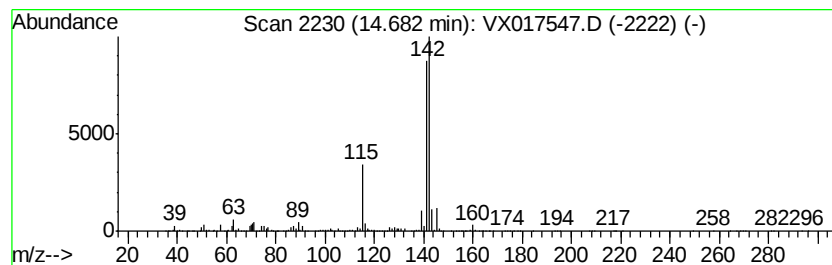
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	63.47 ug/l	3684700	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-dihy...	142	C11H10	004453-90-1	90
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	53



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017547.D
Acq On : 24 Jul 2020 16:53
Operator : JC/SP
Sample : L3413-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
20758

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl Alcohol	2.58	167.8	ug/l	4217470	1	5.64	1256940	50.0
Benzene, 1-ethyl-...	11.42	46.4	ug/l	2693300	4	12.06	2902820	50.0
Benzene, 1,2,3-tr...	12.13	60.4	ug/l	3505180	4	12.06	2902820	50.0
Indane	12.28	77.0	ug/l	4473400	4	12.06	2902820	50.0
Benzene, 2-etheny...	13.21	30.8	ug/l	1787250	4	12.06	2902820	50.0
2,4-Dimethylstyrene	13.33	72.2	ug/l	4191670	4	12.06	2902820	50.0
Naphthalene, 1,2,...	13.46	43.6	ug/l	2533760	4	12.06	2902820	50.0
Naphthalene, 1,2,...	14.27	63.6	ug/l	3694260	4	12.06	2902820	50.0
Naphthalene, 1,2,...	14.52	41.6	ug/l	2414060	4	12.06	2902820	50.0
Naphthalene, 1-me...	14.68	63.5	ug/l	3684700	4	12.06	2902820	50.0