

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX120420\
 Data File : VX019731.D
 Acq On : 04 Dec 2020 20:50
 Operator : JC/MD
 Sample : L4960-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-5

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\82X120420W.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	1.776	107	113	126	rVB	1282068	1746065	22.73%	1.398%
2	1.983	142	147	160	rVB	243765	337607	4.40%	0.270%
3	2.373	204	211	223	rVB	225160	453385	5.90%	0.363%
4	2.825	276	285	288	rBV	884377	1831903	23.85%	1.466%
5	2.867	288	292	304	rVB	1456227	3022554	39.35%	2.419%
6	3.148	328	338	354	rBV	1514141	3440434	44.79%	2.754%
7	3.501	386	396	409	rBV	255052	572722	7.46%	0.458%
8	3.788	436	443	452	rVB	63846	151723	1.98%	0.121%
9	3.898	452	461	474	rVB	52687	137192	1.79%	0.110%
10	4.117	487	497	501	rBV2	49850	145030	1.89%	0.116%
11	4.282	513	524	531	rVV	351238	1067034	13.89%	0.854%
12	4.373	531	539	551	rVV	708511	2037224	26.52%	1.631%
13	4.507	551	561	577	rVB5	37977	158894	2.07%	0.127%
14	5.214	660	677	701	rVB6	80074	425834	5.54%	0.341%
15	5.464	703	718	725	rBV	210693	652137	8.49%	0.522%
16	5.568	725	735	739	rVV5	250829	960382	12.50%	0.769%
17	5.629	740	745	749	rVV	360798	1009694	13.14%	0.808%
18	5.696	749	756	777	rVB	934140	3001179	39.07%	2.402%
19	5.897	777	789	802	rBV3	394350	1237892	16.12%	0.991%
20	6.031	802	811	823	rVB	202722	529270	6.89%	0.424%
21	6.233	825	844	851	rBV2	342059	1191239	15.51%	0.954%
22	6.324	851	859	868	rVV3	826558	2524628	32.87%	2.021%
23	6.428	868	876	887	rVB2	269218	748952	9.75%	0.599%
24	6.671	904	916	928	rVB2	92564	237582	3.09%	0.190%
25	6.830	931	942	948	rBV	511379	1312593	17.09%	1.051%
26	6.909	948	955	967	rVB3	323128	1039959	13.54%	0.832%
27	7.037	968	976	983	rBV	82891	193774	2.52%	0.155%
28	7.123	983	990	999	rVB	78341	176663	2.30%	0.141%
29	7.232	999	1008	1016	rBV	228166	497143	6.47%	0.398%
30	7.434	1029	1041	1052	rBV3	667555	1859479	24.21%	1.488%
31	7.555	1052	1061	1065	rBV2	107510	220230	2.87%	0.176%
32	7.616	1065	1071	1075	rBV	124743	242237	3.15%	0.194%
33	7.714	1081	1087	1097	rVB	180351	359326	4.68%	0.288%
34	7.824	1098	1105	1113	rVB3	108773	222125	2.89%	0.178%

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Integrator: RTE
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35	7.933	1113	1123	1130	rBV5	54171	179695	2.34%	0.144%
36	8.037	1130	1140	1151	rVB	514577	1170540	15.24%	0.937%
37	8.159	1151	1160	1169	rBV2	837255	2184968	28.44%	1.749%
38	8.238	1169	1173	1184	rVB3	179528	392984	5.12%	0.315%
39	8.354	1184	1192	1199	rBV4	136415	276418	3.60%	0.221%
40	8.555	1220	1225	1234	rVB4	153776	322985	4.20%	0.259%
41	8.653	1234	1241	1243	rBV3	211920	394222	5.13%	0.316%
42	8.695	1243	1248	1255	rVB	1095479	1750423	22.79%	1.401%
43	8.964	1285	1292	1297	rBV2	140150	237273	3.09%	0.190%
44	9.019	1297	1301	1306	rVB	102245	155110	2.02%	0.124%
45	9.147	1313	1322	1327	rBV2	204018	388834	5.06%	0.311%
46	9.202	1327	1331	1342	rVB	423350	689922	8.98%	0.552%
47	9.543	1383	1387	1393	rVV4	133400	286262	3.73%	0.229%
48	9.647	1400	1404	1409	rVV2	85077	151191	1.97%	0.121%
49	9.805	1423	1430	1435	rBV3	95127	167099	2.18%	0.134%
50	10.092	1473	1477	1482	rBV	1000681	1356200	17.66%	1.086%
51	10.171	1486	1490	1492	rVV	96543	151672	1.97%	0.121%
52	10.232	1496	1500	1509	rVB	405866	529253	6.89%	0.424%
53	11.000	1621	1626	1634	rBV	1520625	1889486	24.60%	1.512%
54	11.122	1641	1646	1650	rBV	823885	1065332	13.87%	0.853%
55	11.341	1677	1682	1692	rVB	6363624	7681469	100.00%	6.149%
56	11.439	1692	1698	1702	rVV	603151	741059	9.65%	0.593%
57	11.652	1727	1733	1741	rBV	547960	693949	9.03%	0.555%
58	11.786	1751	1755	1761	rVB	327137	412913	5.38%	0.331%
59	11.902	1768	1774	1776	rBV	714524	905576	11.79%	0.725%
60	11.927	1776	1778	1784	rVV	875999	1011763	13.17%	0.810%
61	12.061	1795	1800	1804	rVV	1126094	1402215	18.25%	1.122%
62	12.128	1808	1811	1815	rVB	240310	282567	3.68%	0.226%
63	12.213	1821	1825	1831	rVB	128627	157201	2.05%	0.126%
64	12.286	1831	1837	1843	rBV	4836004	5992307	78.01%	4.797%
65	12.353	1843	1848	1855	rBV2	1678277	2674291	34.81%	2.141%
66	12.445	1858	1863	1867	rVB	642711	788736	10.27%	0.631%
67	12.506	1867	1873	1879	rVB2	1101022	1685773	21.95%	1.349%
68	12.567	1879	1883	1891	rBV	526815	680984	8.87%	0.545%
69	12.652	1891	1897	1900	rBV	6315694	7287201	94.87%	5.833%
70	12.683	1900	1902	1907	rVB	959138	1074377	13.99%	0.860%
71	12.737	1907	1911	1919	rBV2	2589880	3717684	48.40%	2.976%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON Filtering: 5
 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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72	12.884	1930	1935	1940	rVB3	377019	535443	6.97%	0.429%
73	12.963	1940	1948	1951	rBV	4935261	6108363	79.52%	4.889%
74	13.000	1951	1954	1960	rVV	1780772	2102253	27.37%	1.683%
75	13.061	1960	1964	1973	rVV2	724532	1011508	13.17%	0.810%
76	13.152	1973	1979	1980	rVV2	234415	359463	4.68%	0.288%
77	13.176	1980	1983	1985	rVV	437758	583339	7.59%	0.467%
78	13.207	1985	1988	1998	rVV	2840212	3487111	45.40%	2.791%
79	13.298	1998	2003	2004	rVV	359815	427296	5.56%	0.342%
80	13.329	2004	2008	2014	rVV2	4865836	6817847	88.76%	5.457%
81	13.384	2014	2017	2018	rVV2	269772	337548	4.39%	0.270%
82	13.408	2018	2021	2026	rVV	718069	905701	11.79%	0.725%
83	13.463	2026	2030	2037	rVB	518968	662820	8.63%	0.531%
84	13.542	2037	2043	2046	rBV2	535967	744015	9.69%	0.596%
85	13.585	2046	2050	2052	rVV	1002706	1349774	17.57%	1.080%
86	13.621	2052	2056	2061	rVV3	1623400	2760393	35.94%	2.210%
87	13.682	2062	2066	2067	rVV	516210	754765	9.83%	0.604%
88	13.701	2067	2069	2076	rVB2	1139553	1832896	23.86%	1.467%
89	13.817	2084	2088	2097	rVB	286412	440616	5.74%	0.353%
90	13.896	2097	2101	2105	rVB	156631	217524	2.83%	0.174%
91	13.963	2105	2112	2116	rBV2	568939	785025	10.22%	0.628%
92	14.012	2116	2120	2124	rBV	465277	565459	7.36%	0.453%
93	14.115	2132	2137	2143	rBV	1047101	1255500	16.34%	1.005%
94	14.255	2155	2160	2165	rBV	717010	1065523	13.87%	0.853%
95	14.359	2173	2177	2182	rBV	515823	639926	8.33%	0.512%
96	14.414	2182	2186	2190	rVB	488625	619215	8.06%	0.496%
97	14.524	2198	2204	2208	rBV2	161169	247267	3.22%	0.198%
98	14.676	2216	2229	2234	rBV	1293263	1780184	23.18%	1.425%
99	14.822	2247	2253	2257	rBV	1323283	1696179	22.08%	1.358%
100	15.664	2384	2391	2395	rBV	96214	162630	2.12%	0.130%

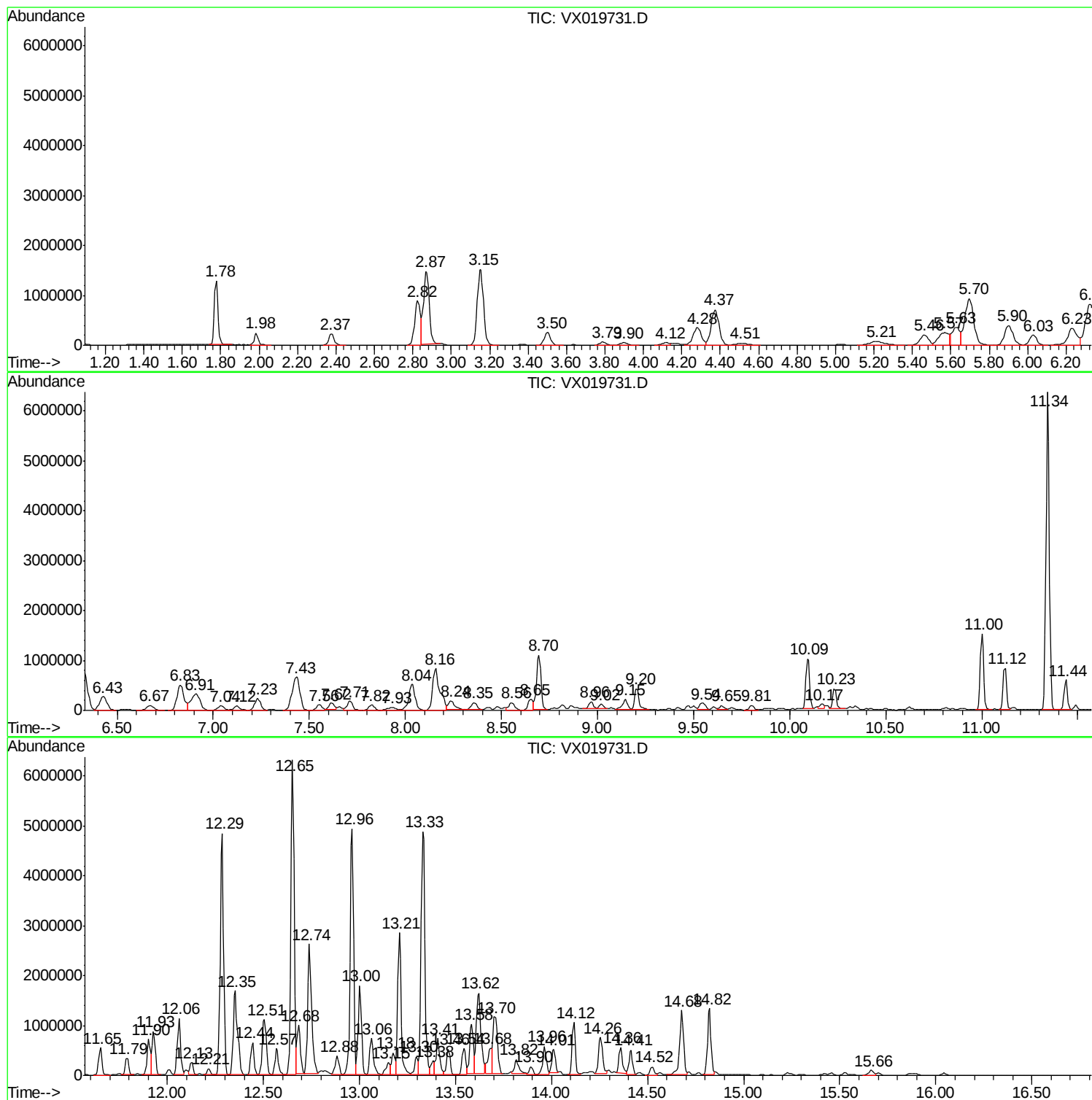
Sum of corrected areas: 124929602

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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



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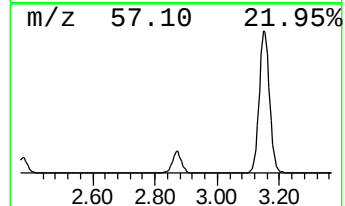
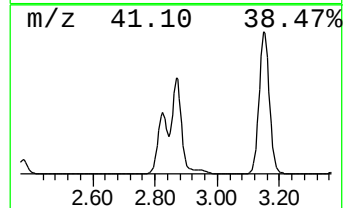
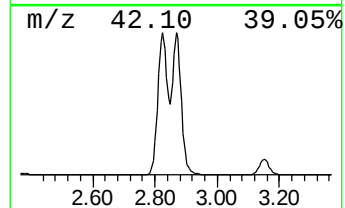
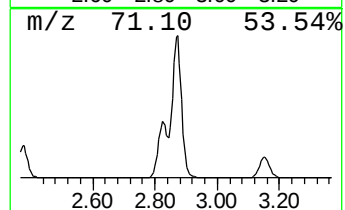
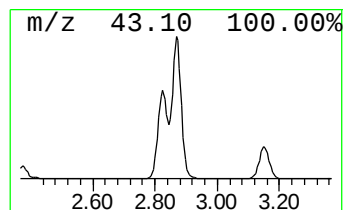
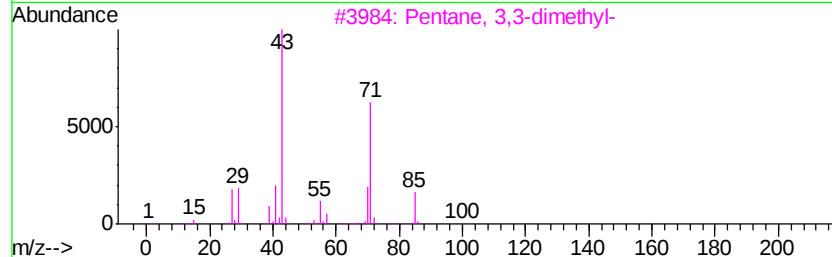
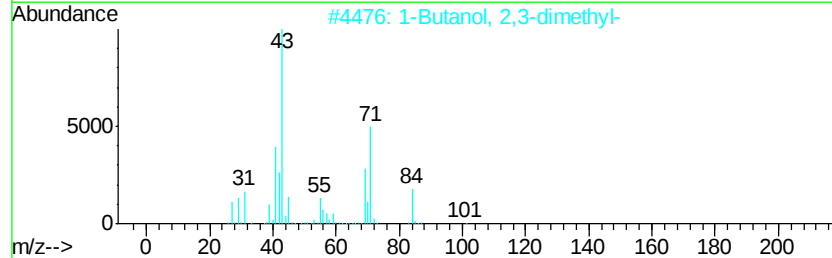
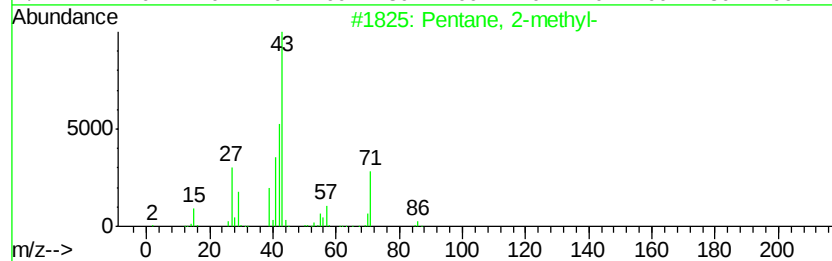
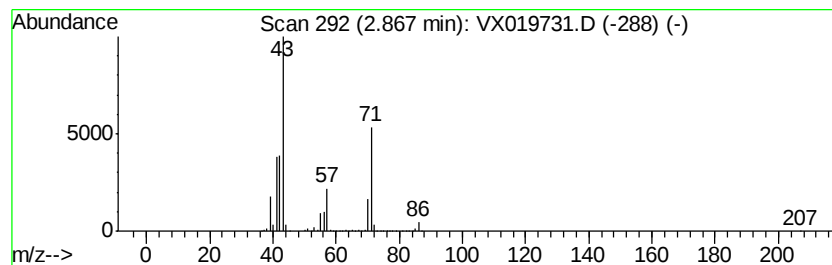
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	149.68 ug/l	3022550	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2-methyl-	86	C6H14	000107-83-5	83
2		1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
3		Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
4		Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	36
5		3-Pyrrolidinol	87	C4H9NO	040499-83-0	33



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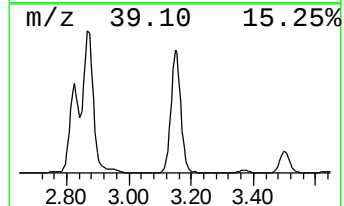
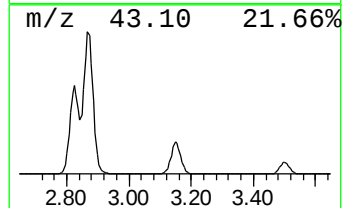
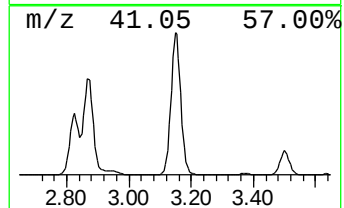
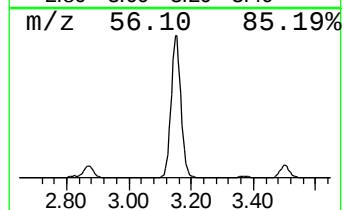
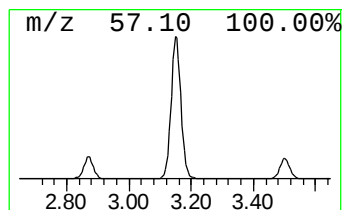
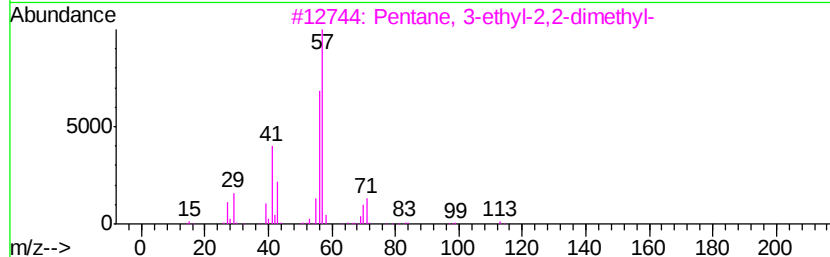
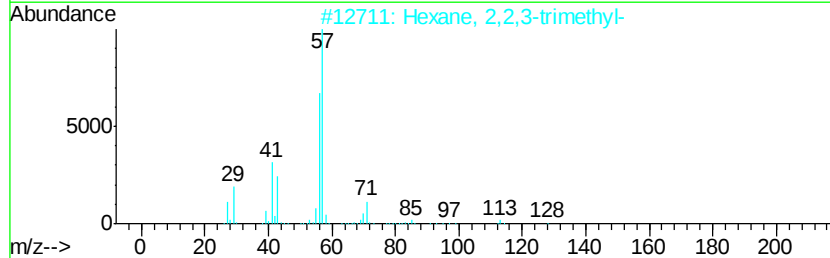
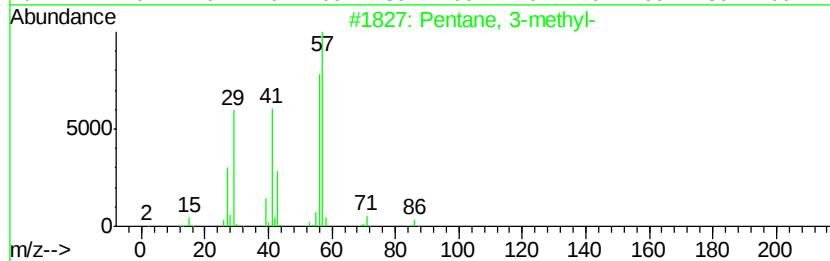
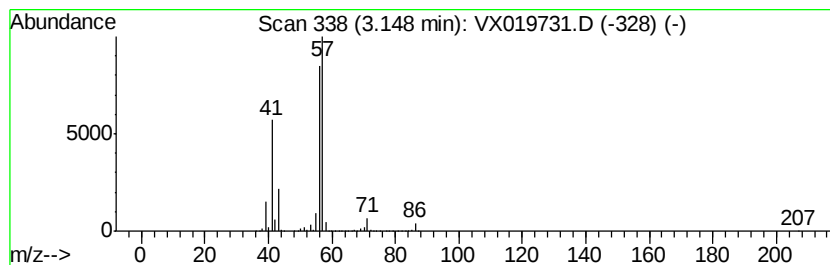
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.15	170.37 ug/l	3440430	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4		n-Hexane	86	C6H14	000110-54-3	47
5		Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45



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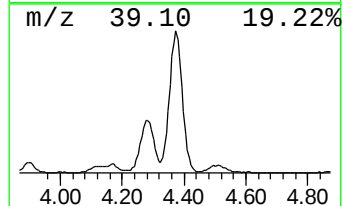
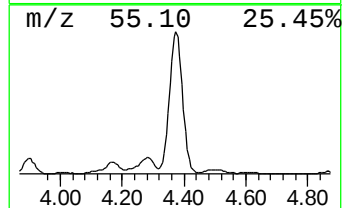
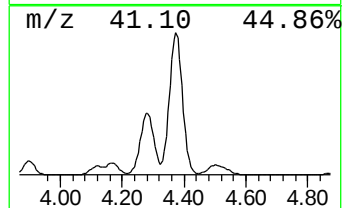
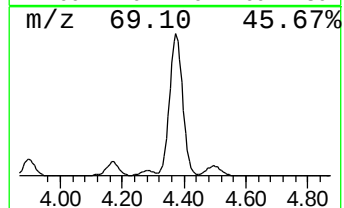
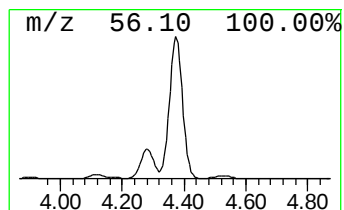
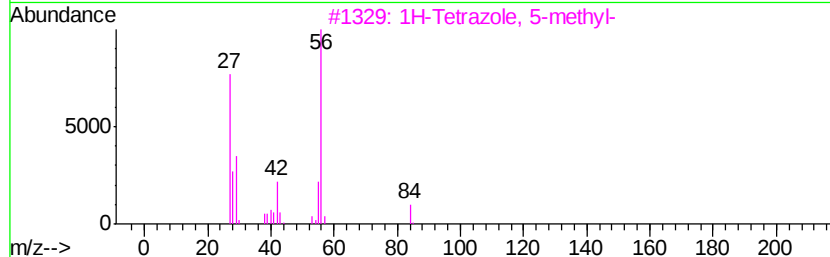
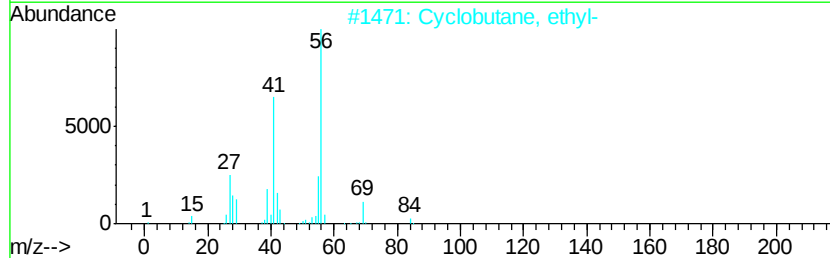
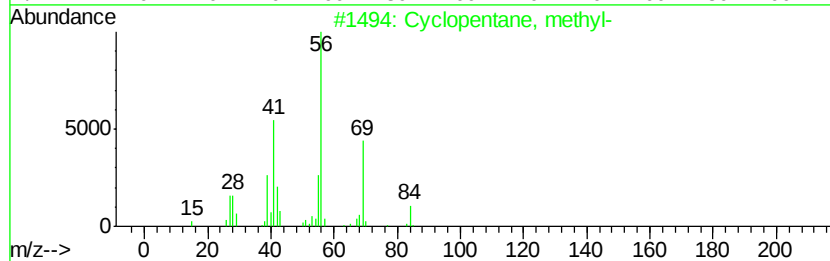
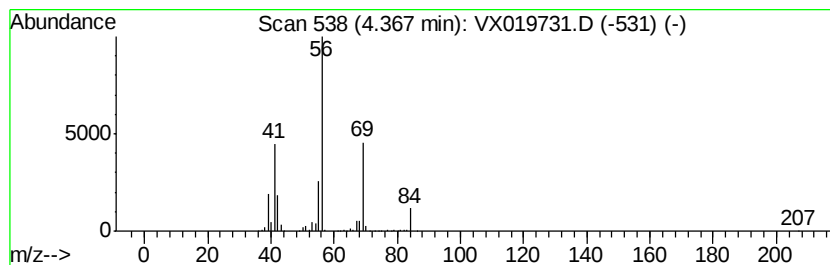
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Quant Title : SW846 8260

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

Peak Number 3 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	100.88 ug/l	2037220	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4		Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	43
5		2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	36



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

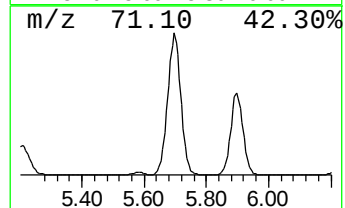
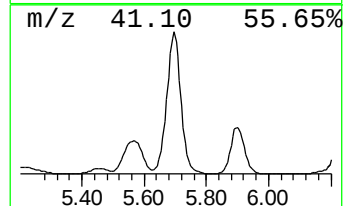
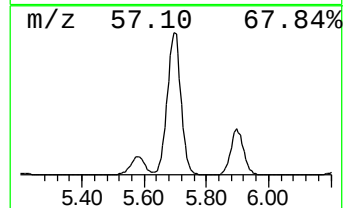
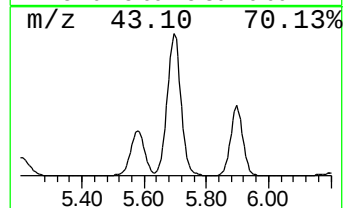
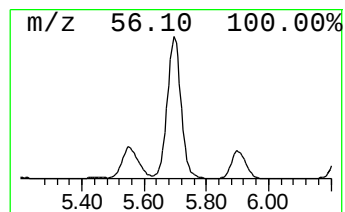
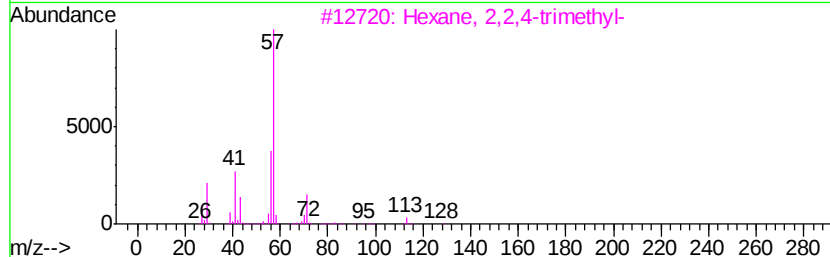
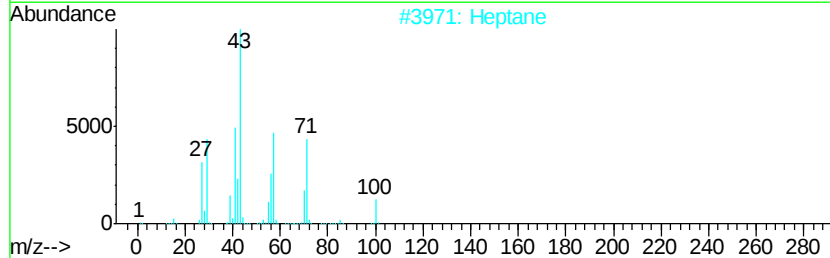
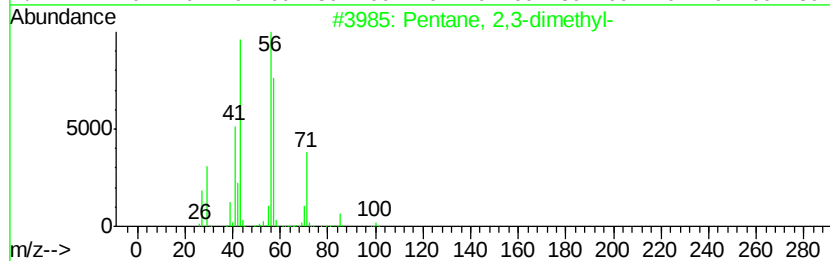
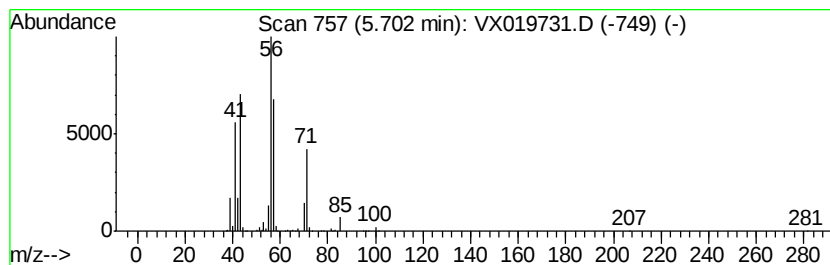
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Quant Title : SW846 8260

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

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	148.62 ug/l	3001180	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2		Heptane	100	C7H16	000142-82-5	47
3		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	39
4		1-Hexene, 5-methyl-	98	C7H14	003524-73-0	38
5		Aziridine, 2-methyl-	57	C3H7N	000075-55-8	37



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

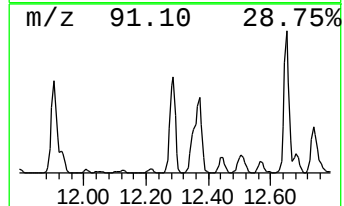
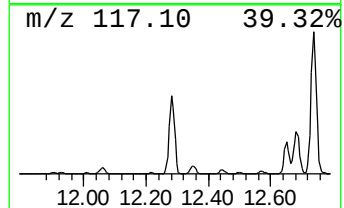
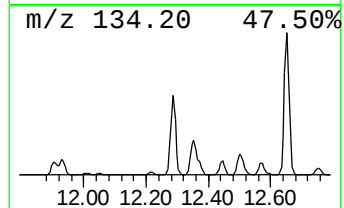
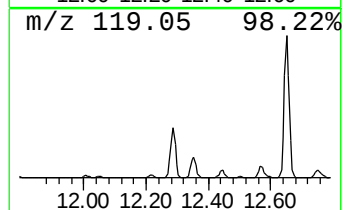
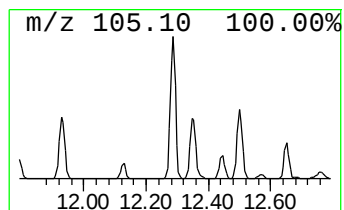
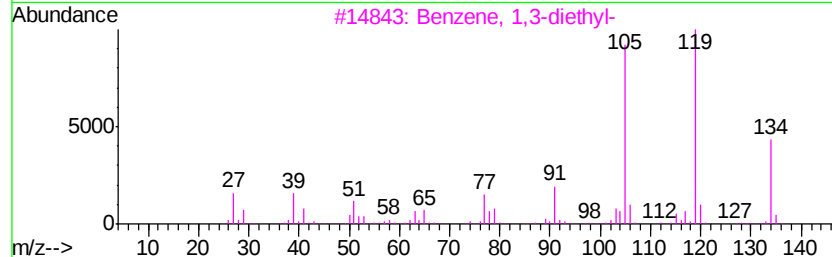
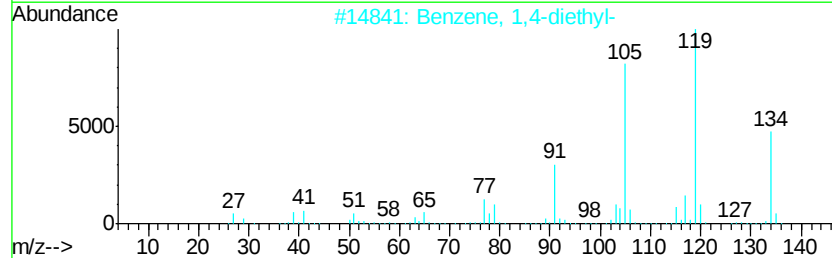
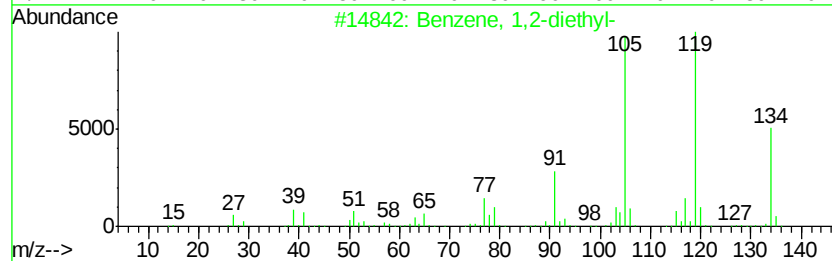
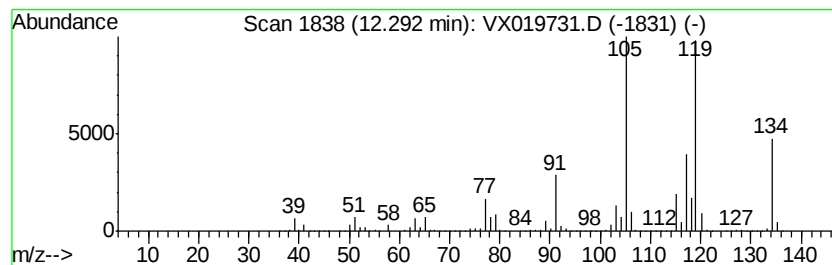
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	213.67 ug/l	5992310	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	70
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	62
5		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

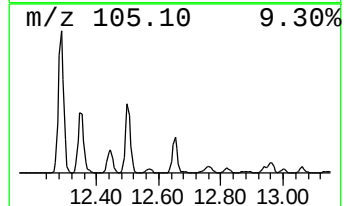
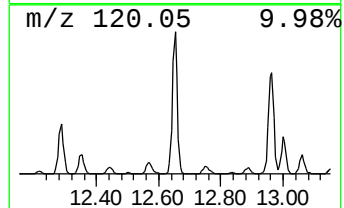
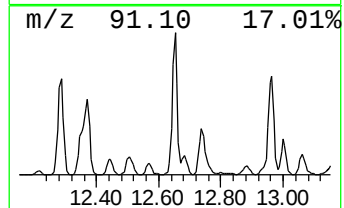
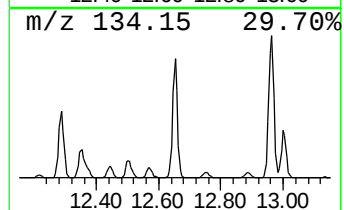
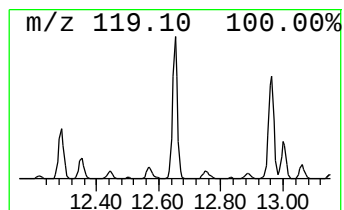
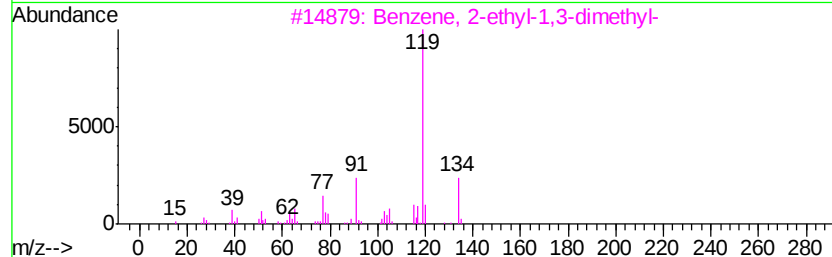
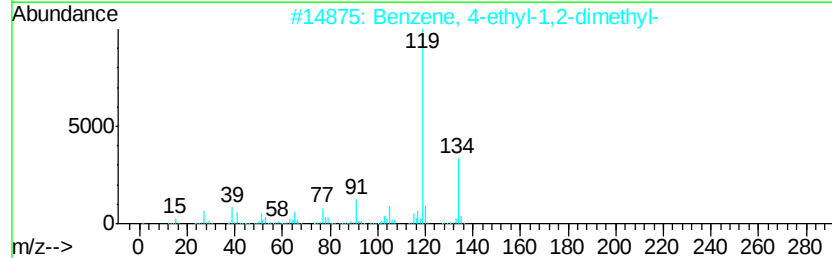
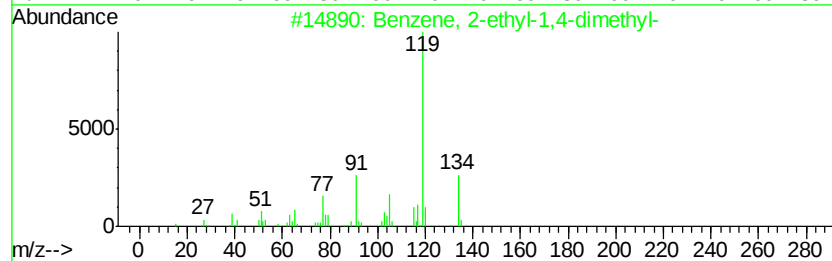
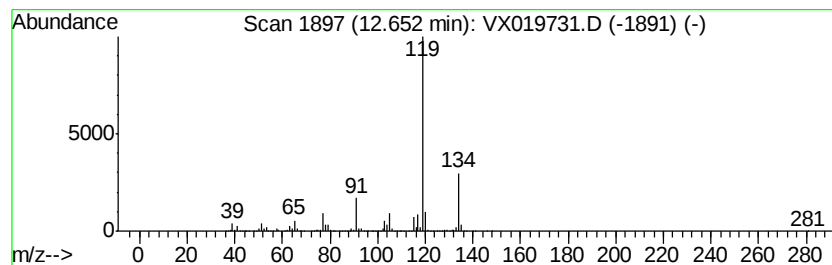
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	259.85 ug/l	7287200	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-5

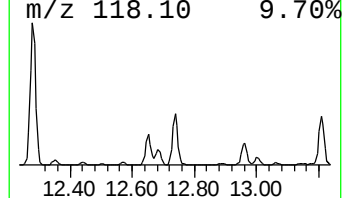
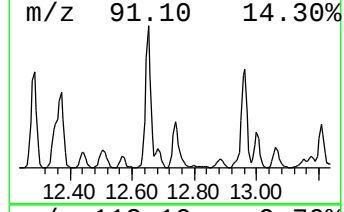
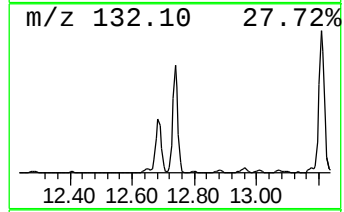
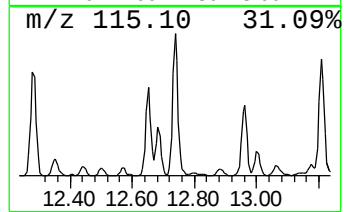
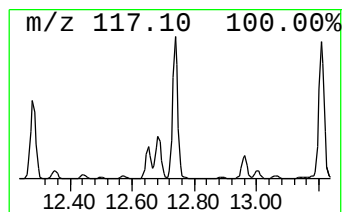
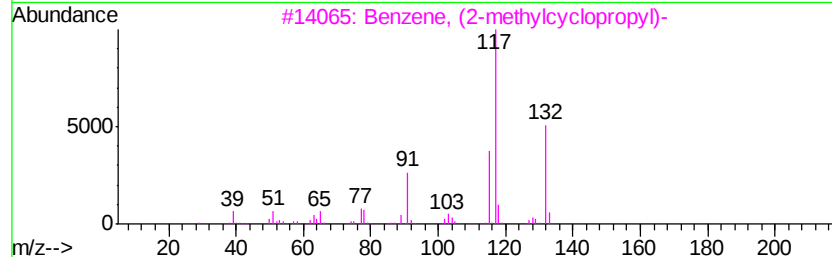
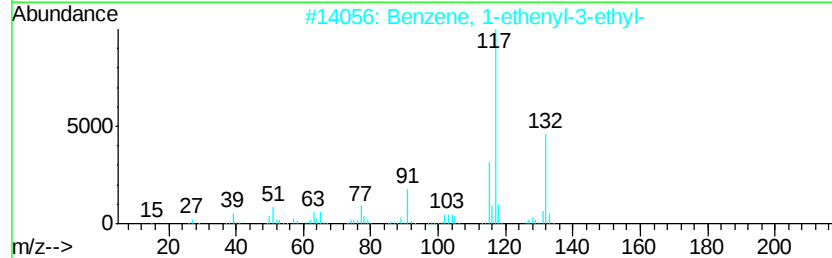
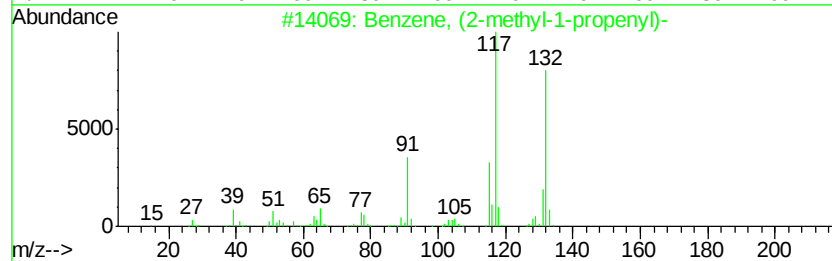
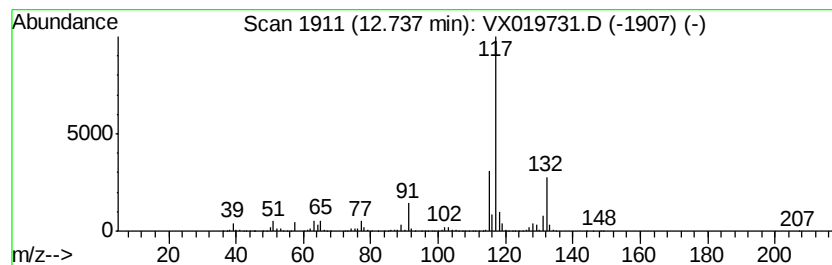
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	132.57 ug/l	3717680	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5		Indan, 1-methyl-	132	C10H12	000767-58-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-5

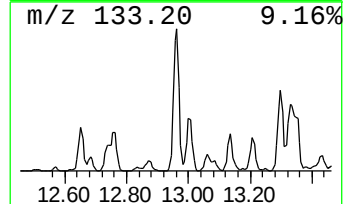
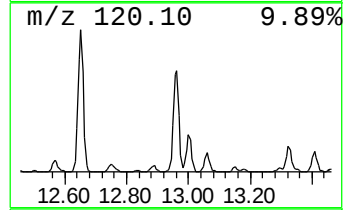
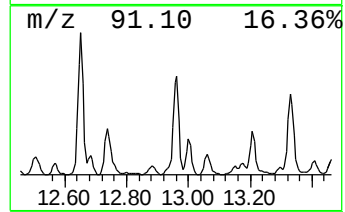
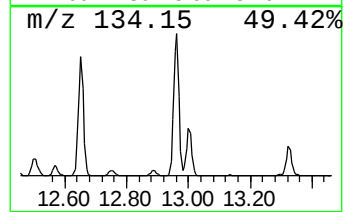
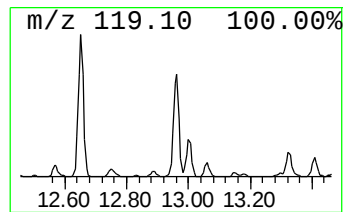
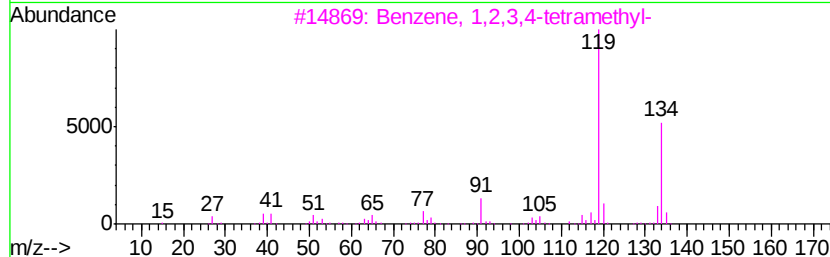
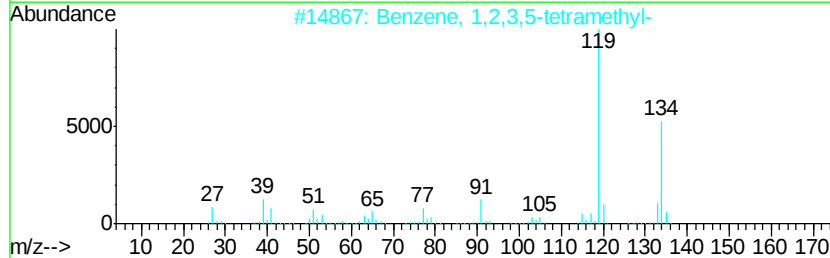
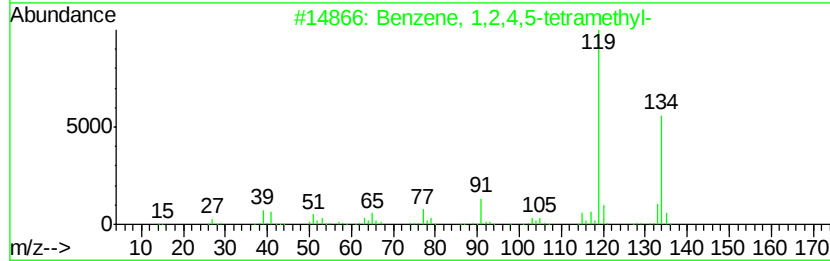
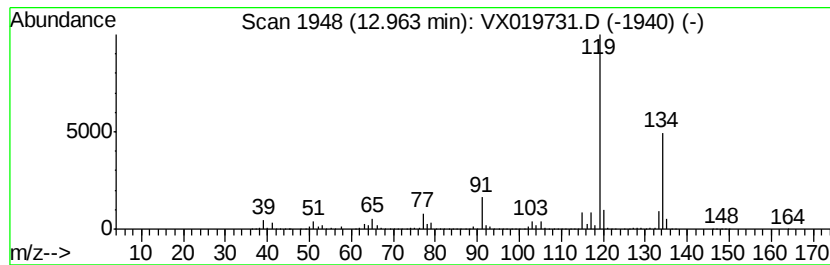
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	217.81 ug/l	6108360	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-5

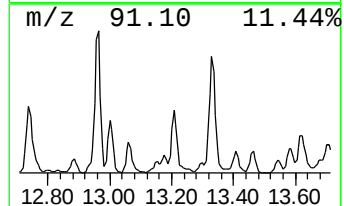
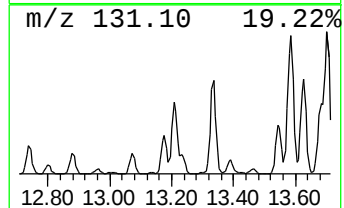
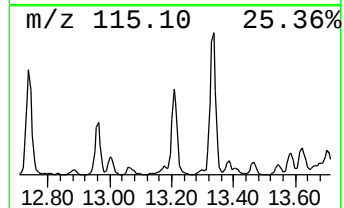
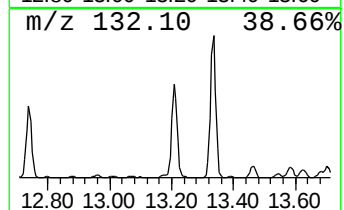
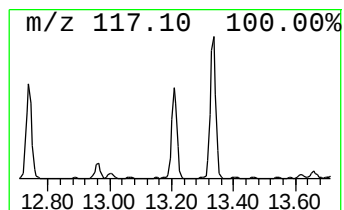
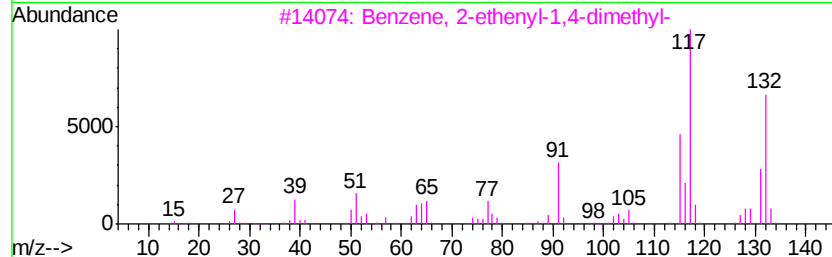
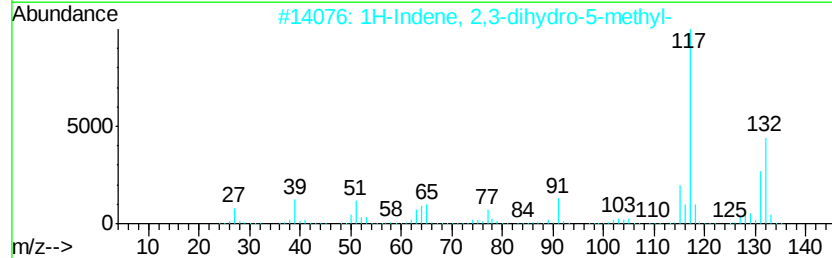
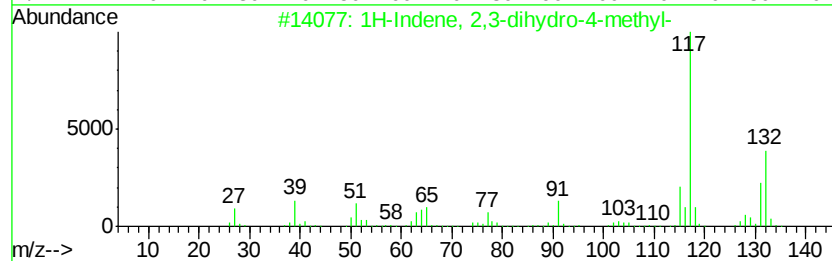
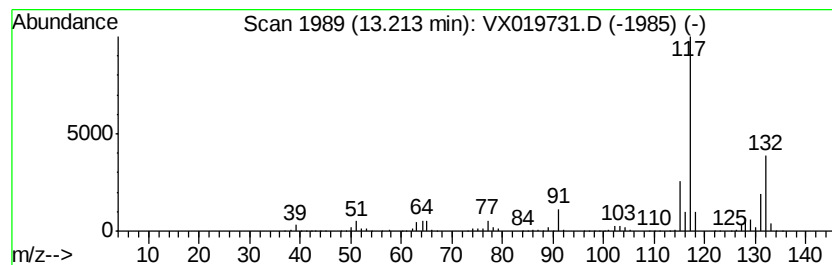
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	124.34 ug/l	3487110	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
MW-5

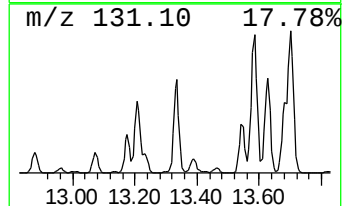
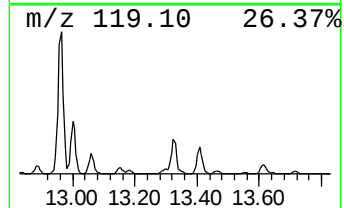
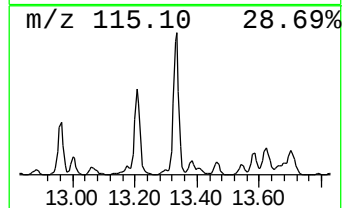
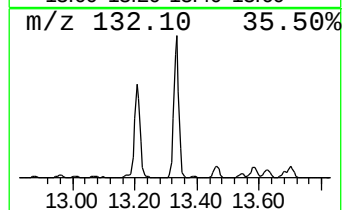
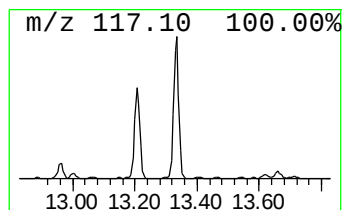
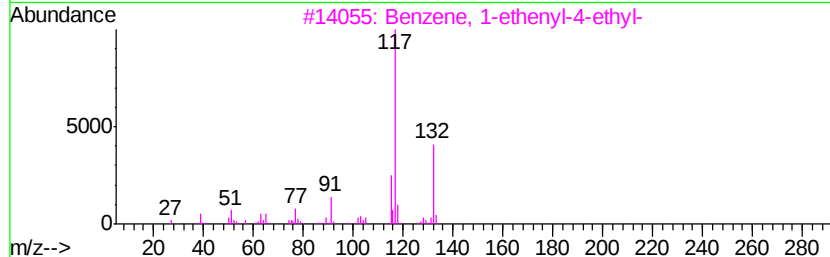
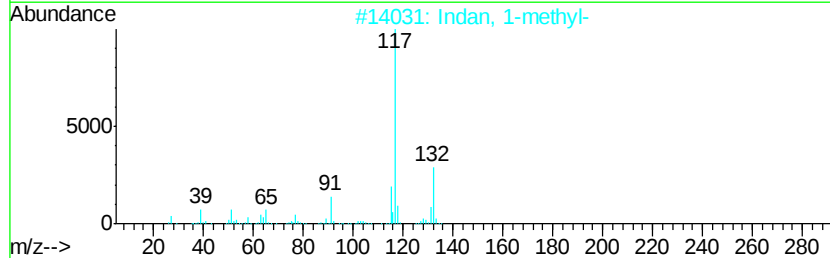
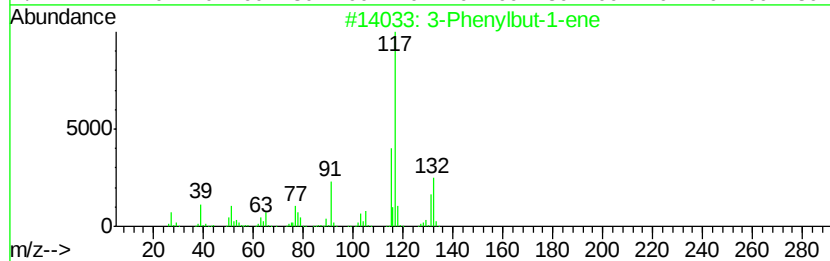
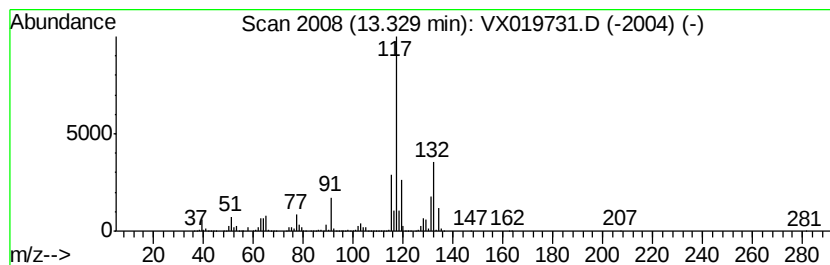
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Quant Title : SW846 8260

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

Peak Number 10 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	243.11 ug/l	6817850	1,4-Dichlorobenzene-d4	12.06

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
2			Indan, 1-methyl-	132	C10H12	000767-58-8	81
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	74
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX120420\
Data File : VX019731.D
Acq On : 04 Dec 2020 20:50
Operator : JC/MD
Sample : L4960-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.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
Pentane, 2-methyl-	2.87	149.7	ug/l	3022550	1	5.63	1009690	50.0
Pentane, 3-methyl-	3.15	170.4	ug/l	3440430	1	5.63	1009690	50.0
Cyclopentane, met...	4.37	100.9	ug/l	2037220	1	5.63	1009690	50.0
Pentane, 2,3-dime...	5.70	148.6	ug/l	3001180	1	5.63	1009690	50.0
Benzene, 1,2-diet...	12.29	213.7	ug/l	5992310	4	12.06	1402220	50.0
Benzene, 2-ethyl-...	12.65	259.9	ug/l	7287200	4	12.06	1402220	50.0
Benzene, (2-methy...	12.74	132.6	ug/l	3717680	4	12.06	1402220	50.0
Benzene, 1,2,4,5-...	12.96	217.8	ug/l	6108360	4	12.06	1402220	50.0
1H-Indene, 2,3-di...	13.21	124.3	ug/l	3487110	4	12.06	1402220	50.0
3-Phenylbut-1-ene	13.33	243.1	ug/l	6817850	4	12.06	1402220	50.0