

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX010320\
 Data File : VX014400.D
 Acq On : 03 Jan 2020 15:04
 Operator : JC/SP
 Sample : K6476-01
 Misc : 6.96g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WP022SB454_191227-12-14

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\82X121319W.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	3.496	418	428	438	rBV	313230	711583	1.28%	0.164%
2	5.490	744	755	761	rBV	246578	755449	1.36%	0.174%
3	5.588	761	771	776	rVV4	337300	1092473	1.97%	0.252%
4	5.648	776	781	788	rVB	317913	791315	1.43%	0.182%
5	5.911	811	824	838	rBV	530004	1593477	2.88%	0.367%
6	6.057	838	848	854	rVV	285836	787180	1.42%	0.181%
7	6.331	885	893	903	rVV	505720	1554257	2.81%	0.358%
8	6.691	941	952	963	rBV	1123579	2668899	4.82%	0.615%
9	6.849	969	978	998	rVB	721137	1842080	3.33%	0.424%
10	7.453	1066	1077	1087	rBV	1385824	3208271	5.79%	0.739%
11	8.044	1162	1174	1185	rBV3	310638	893242	1.61%	0.206%
12	8.331	1216	1221	1225	rBV	858957	1428026	2.58%	0.329%
13	8.496	1243	1248	1258	rBV	768207	1336916	2.41%	0.308%
14	8.660	1268	1275	1279	rBV3	800796	1685160	3.04%	0.388%
15	8.709	1279	1283	1289	rVB	1609039	2757811	4.98%	0.635%
16	8.782	1289	1295	1300	rBV	1084578	1761200	3.18%	0.406%
17	8.965	1320	1325	1331	rBV	2193917	3119841	5.63%	0.719%
18	9.032	1331	1336	1344	rVB2	434068	746013	1.35%	0.172%
19	9.440	1393	1403	1412	rBV	515984	856368	1.55%	0.197%
20	9.556	1416	1422	1428	rBV3	513128	1162675	2.10%	0.268%
21	9.617	1428	1432	1436	rVB	866067	1182616	2.14%	0.272%
22	9.672	1436	1441	1445	rBV	787396	1188416	2.15%	0.274%
23	9.886	1468	1476	1479	rBV3	632016	1497719	2.70%	0.345%
24	9.959	1483	1488	1493	rVB	2011137	3023483	5.46%	0.696%
25	10.062	1498	1505	1509	rBV	2061031	2891261	5.22%	0.666%
26	10.135	1509	1517	1529	rVB	22532852	34488431	62.28%	7.945%
27	10.245	1529	1535	1540	rBV2	791424	1402635	2.53%	0.323%
28	10.355	1540	1553	1558	rBV2	2882738	6971241	12.59%	1.606%
29	10.410	1558	1562	1573	rVB	12012680	15454268	27.91%	3.560%
30	10.507	1573	1578	1584	rBV2	531846	796529	1.44%	0.183%
31	10.635	1593	1599	1606	rVV2	2140584	3786791	6.84%	0.872%
32	10.696	1606	1609	1611	rVV	1036055	1507477	2.72%	0.347%
33	10.721	1611	1613	1617	rVV	1411007	1841648	3.33%	0.424%
34	10.763	1617	1620	1623	rVB2	621759	779595	1.41%	0.180%

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35	10.830	1623	1631	1635	rBV2	4992322	7384683	13.34%	1.701%
36	10.910	1635	1644	1647	rVV2	4079693	7274661	13.14%	1.676%
37	10.946	1647	1650	1655	rVV	2441555	3326587	6.01%	0.766%
38	11.001	1655	1659	1664	rVV3	724266	1374486	2.48%	0.317%
39	11.056	1664	1668	1669	rVV	896873	1210464	2.19%	0.279%
40	11.080	1669	1672	1675	rVV	2321421	3278920	5.92%	0.755%
41	11.135	1675	1681	1683	rVV2	5521163	8728031	15.76%	2.011%
42	11.160	1683	1685	1691	rVB2	3737071	4388470	7.92%	1.011%
43	11.239	1691	1698	1701	rBV	4256652	6178785	11.16%	1.423%
44	11.269	1701	1703	1712	rVV3	2613987	3658533	6.61%	0.843%
45	11.355	1713	1717	1720	rVV	933573	1303352	2.35%	0.300%
46	11.391	1720	1723	1726	rVV2	1047498	1704031	3.08%	0.393%
47	11.434	1726	1730	1735	rVV2	3265943	6512608	11.76%	1.500%
48	11.483	1735	1738	1740	rVV	3555070	5120194	9.25%	1.179%
49	11.532	1741	1746	1750	rVV	21909380	29013229	52.39%	6.683%
50	11.574	1750	1753	1758	rVV5	479168	820095	1.48%	0.189%
51	11.623	1758	1761	1764	rVV	535498	799611	1.44%	0.184%
52	11.666	1764	1768	1775	rVV5	1899651	4461812	8.06%	1.028%
53	11.733	1775	1779	1781	rVV2	1146548	1536797	2.78%	0.354%
54	11.763	1781	1784	1787	rVV	3631802	4401203	7.95%	1.014%
55	11.806	1787	1791	1800	rVB	5953200	9730121	17.57%	2.241%
56	11.885	1800	1804	1806	rBV2	902083	1183402	2.14%	0.273%
57	11.916	1806	1809	1810	rVV	1137145	1375099	2.48%	0.317%
58	11.940	1810	1813	1818	rVV3	2126329	3619444	6.54%	0.834%
59	11.995	1818	1822	1824	rVV	4055149	5409632	9.77%	1.246%
60	12.025	1824	1827	1830	rVV	6012233	7642395	13.80%	1.760%
61	12.092	1830	1838	1842	rVV	19833939	29184841	52.70%	6.723%
62	12.147	1842	1847	1852	rVV3	3281031	6694246	12.09%	1.542%
63	12.220	1852	1859	1867	rVB3	2321203	4383322	7.92%	1.010%
64	12.318	1867	1875	1879	rBV2	2592933	4965510	8.97%	1.144%
65	12.391	1879	1887	1892	rVV	38104004	55376564	100.00%	12.756%
66	12.470	1892	1900	1904	rVV	11051492	14400740	26.01%	3.317%
67	12.519	1904	1908	1914	rVB2	2175413	3602322	6.51%	0.830%
68	12.605	1914	1922	1928	rBV3	2583622	6236228	11.26%	1.437%
69	12.665	1928	1932	1939	rVB	1958250	2905163	5.25%	0.669%
70	12.769	1944	1949	1954	rVV4	2313785	4450463	8.04%	1.025%
71	12.848	1954	1962	1963	rVV5	1068787	2255818	4.07%	0.520%

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72	12.873	1963	1966	1968	rVV	1605137	2330248	4.21%	0.537%
73	12.897	1968	1970	1973	rVV2	1077336	1259227	2.27%	0.290%
74	12.940	1973	1977	1980	rVV3	1822329	2919363	5.27%	0.672%
75	12.970	1980	1982	1986	rVV2	1880865	2518408	4.55%	0.580%
76	13.013	1986	1989	1996	rVV2	2582245	4472093	8.08%	1.030%
77	13.080	1996	2000	2001	rVV2	928716	1247752	2.25%	0.287%
78	13.123	2005	2007	2009	rVV	934776	940518	1.70%	0.217%
79	13.147	2009	2011	2016	rVB2	527222	691220	1.25%	0.159%
80	13.208	2016	2021	2026	rVB3	1246541	1899394	3.43%	0.438%
81	13.263	2026	2030	2033	rBV	749413	834501	1.51%	0.192%
82	13.312	2033	2038	2040	rBV3	1228866	2092059	3.78%	0.482%
83	13.342	2040	2043	2048	rVV2	2970438	4392444	7.93%	1.012%
84	13.476	2059	2065	2075	rVB7	1356601	2946823	5.32%	0.679%
85	13.641	2087	2092	2097	rBV3	3524436	5320243	9.61%	1.226%
86	13.726	2102	2106	2110	rVB3	516601	783226	1.41%	0.180%
87	13.830	2119	2123	2130	rVB	3332898	4518246	8.16%	1.041%
88	13.982	2140	2148	2151	rBV4	493106	1055481	1.91%	0.243%
89	14.269	2190	2195	2201	rBV4	529562	958542	1.73%	0.221%
90	14.354	2206	2209	2217	rVB2	363326	760812	1.37%	0.175%
91	14.537	2235	2239	2247	rVB5	347836	679997	1.23%	0.157%
92	14.690	2257	2264	2268	rBV	3498927	4551828	8.22%	1.049%
93	14.836	2281	2288	2296	rVB	2161585	3081364	5.56%	0.710%
94	15.147	2335	2339	2343	rVB2	1144515	1513991	2.73%	0.349%
95	15.281	2350	2361	2364	rBV4	535556	1455081	2.63%	0.335%
96	15.439	2380	2387	2390	rBV3	572433	833007	1.50%	0.192%
97	15.543	2396	2404	2412	rBV	1330493	2187261	3.95%	0.504%
98	15.683	2421	2427	2430	rBV	1266561	2011281	3.63%	0.463%
99	15.720	2430	2433	2440	rVB2	980507	1460337	2.64%	0.336%
100	15.909	2457	2464	2474	rVB3	310364	943207	1.70%	0.217%

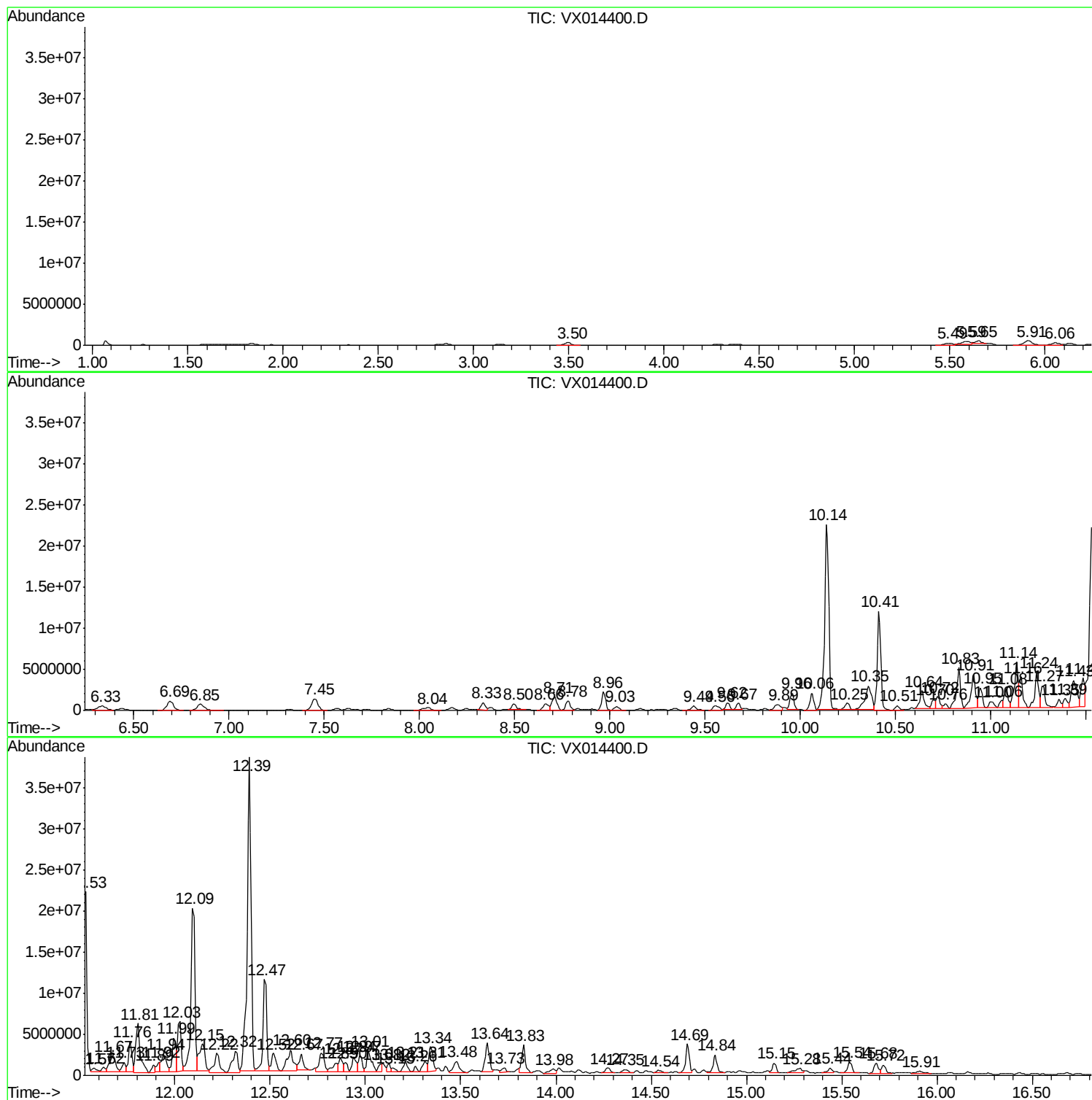
Sum of corrected areas: 434114162

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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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



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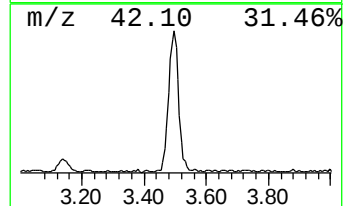
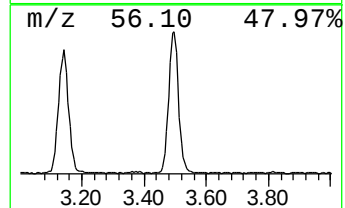
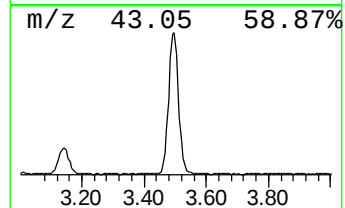
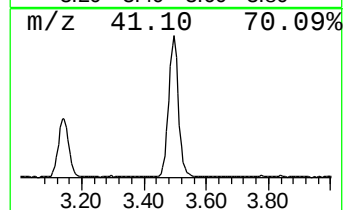
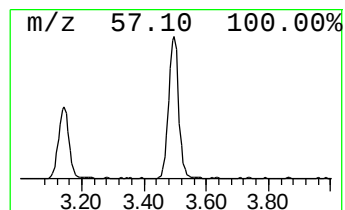
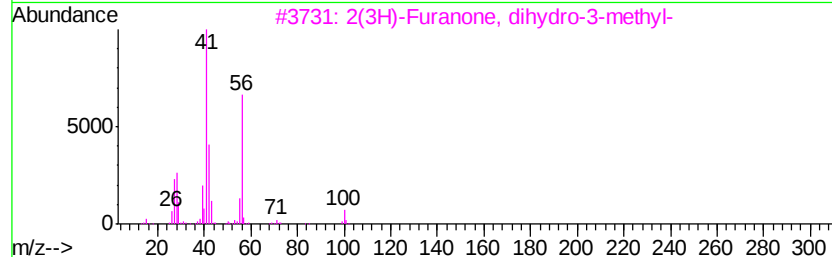
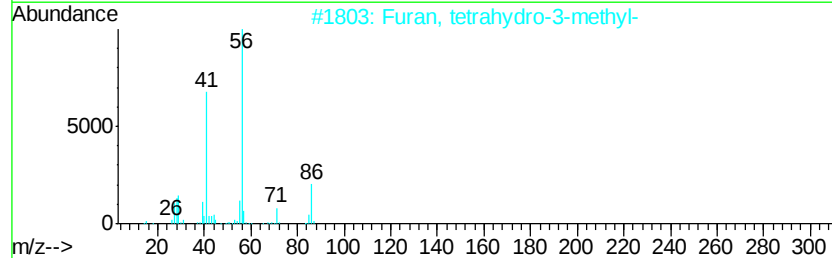
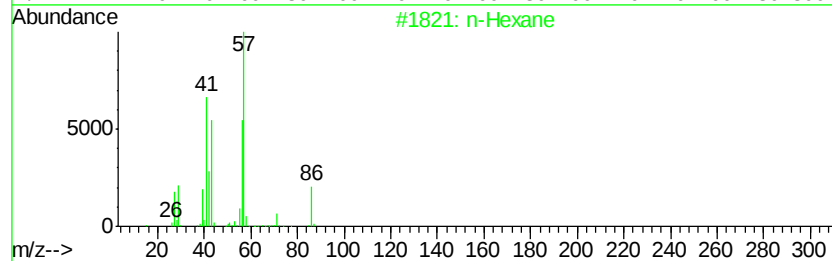
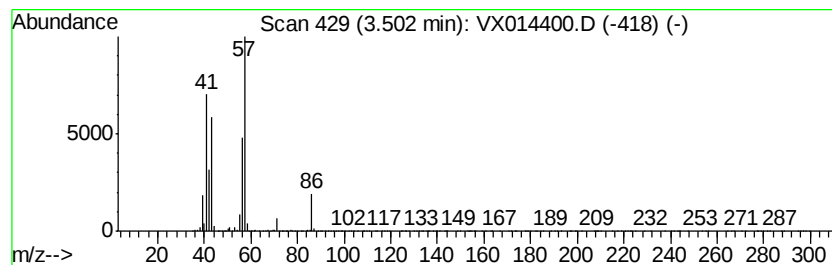
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 1 n-Hexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.50	44.96 ug/l	711583	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexane	86	C6H14	000110-54-3	94
2		Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	53
3		2(3H)-Furanone, dihydro-3-methyl-	100	C5H8O2	001679-47-6	50
4		Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43
5		3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	38



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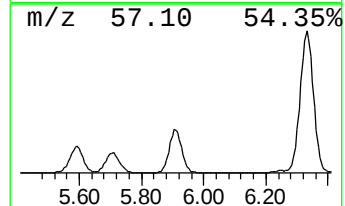
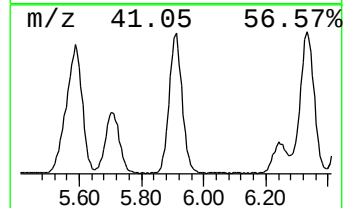
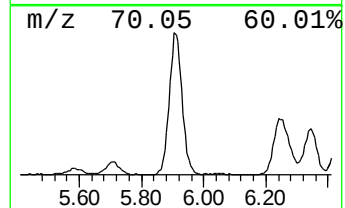
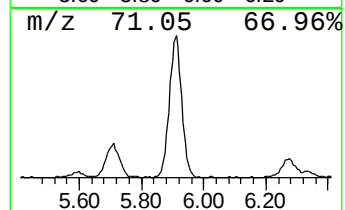
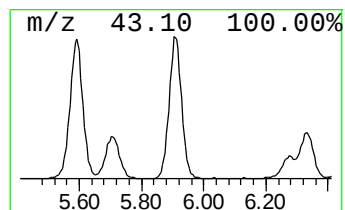
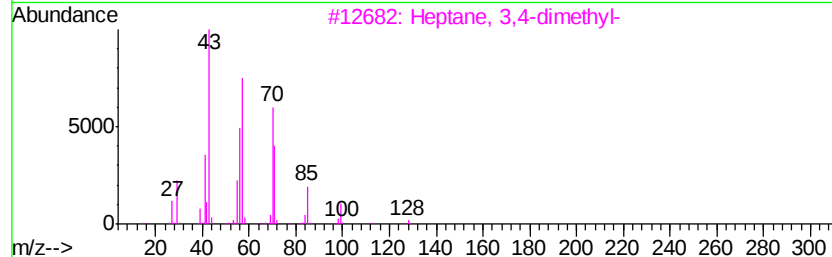
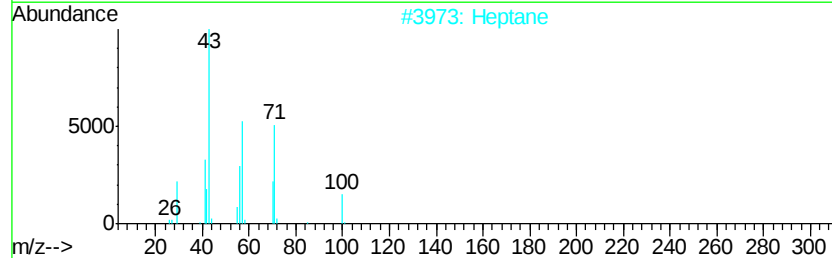
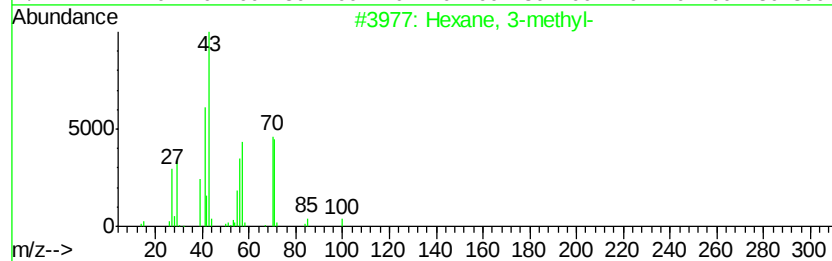
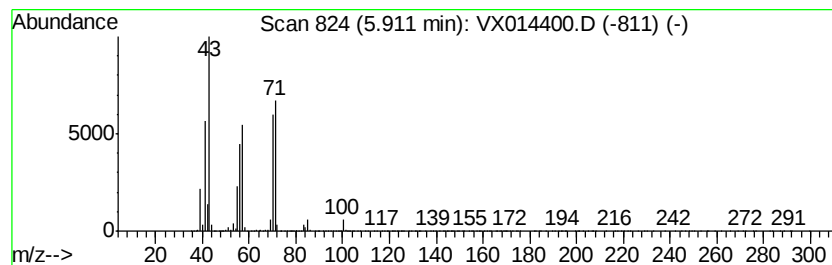
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Quant Title : SW846 8260

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

Peak Number 2 Hexane, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.91	100.69 ug/l	1593480	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2		Heptane	100	C7H16	000142-82-5	80
3		Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	59
4		Heptane, 4-methyl-	114	C8H18	000589-53-7	59
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	53



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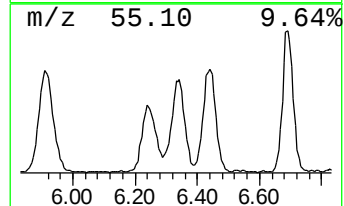
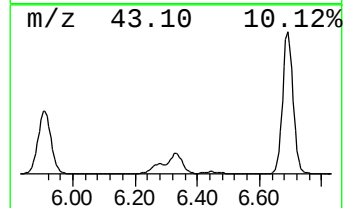
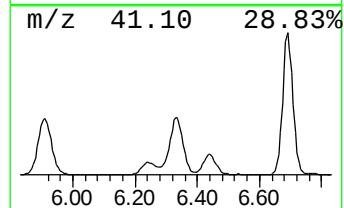
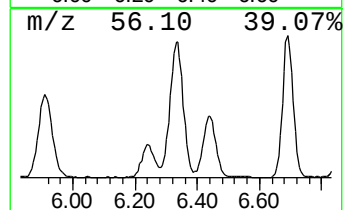
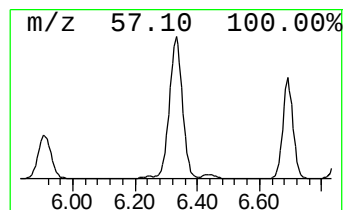
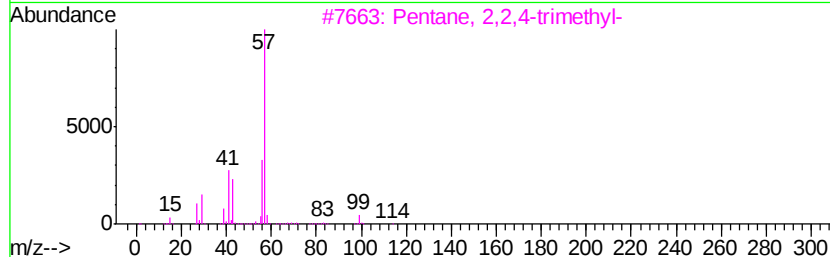
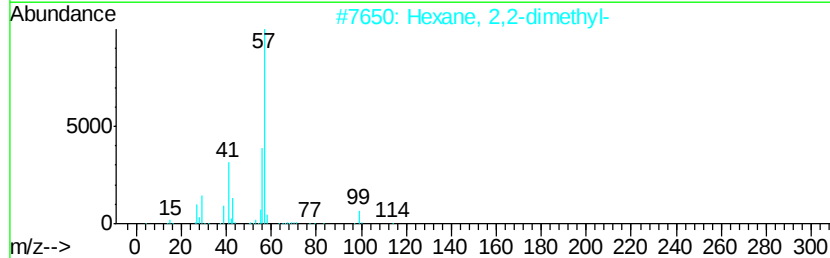
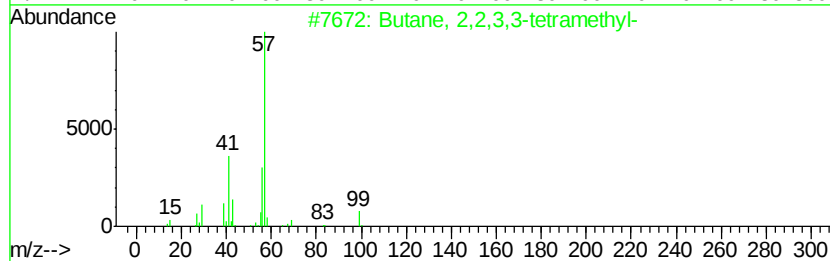
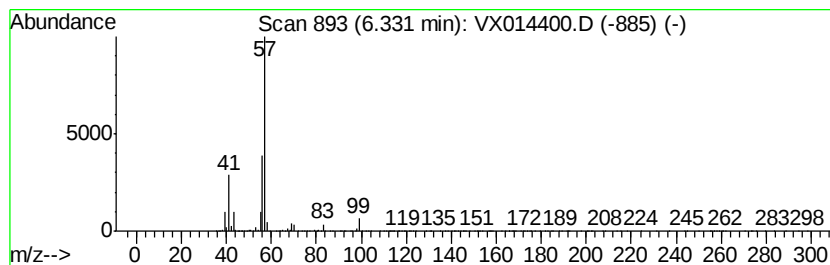
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Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,2,3,3-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	42.19 ug/l	1554260	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
4		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WP022SB454_191227-12-14

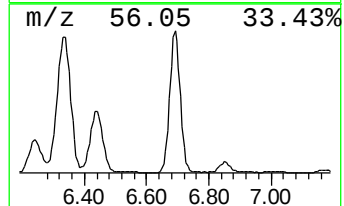
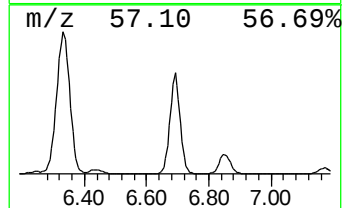
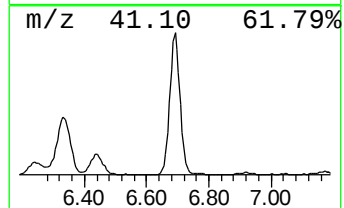
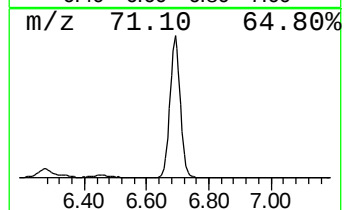
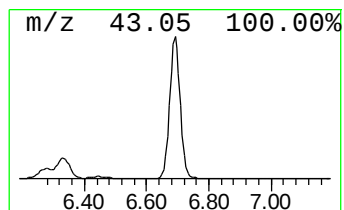
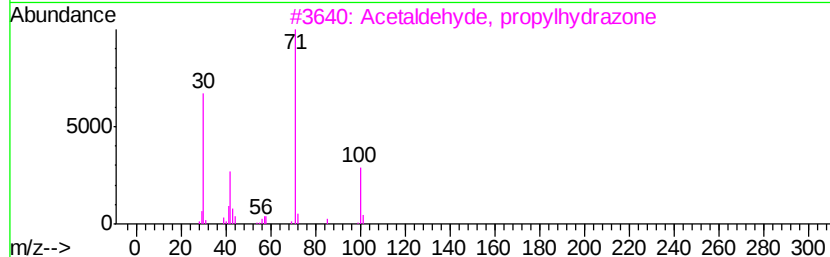
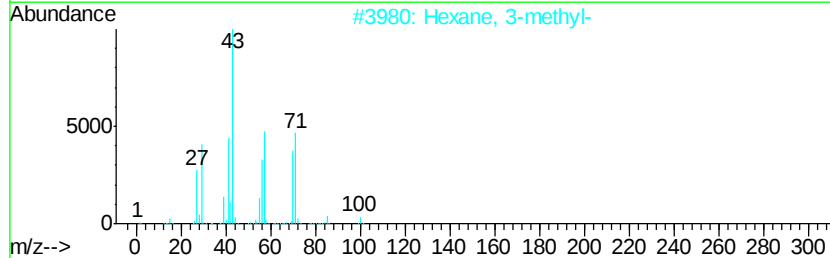
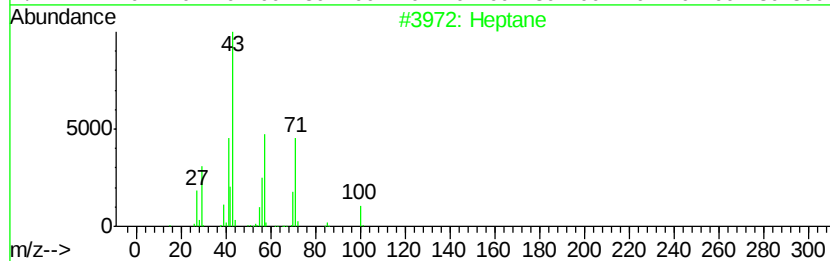
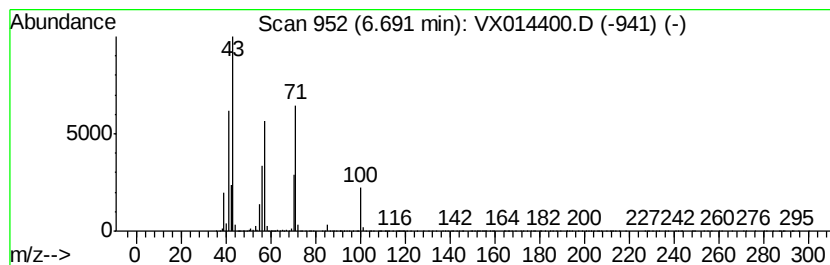
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Quant Title : SW846 8260

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

Peak Number 4 Heptane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	72.44 ug/l	2668900	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane	100	C7H16	000142-82-5	91
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	47
3		Acetaldehyde, propylhydrazine	100	C5H12N2	007422-88-0	47
4		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	40
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	37



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WP022SB454_191227-12-14

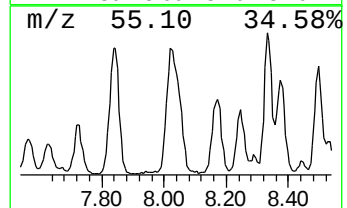
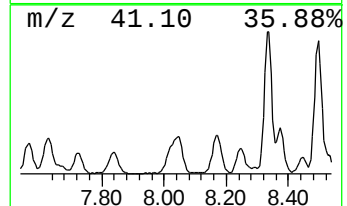
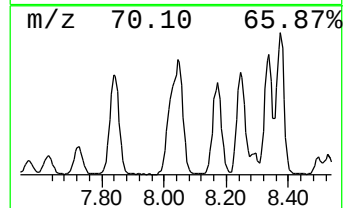
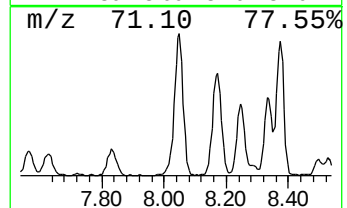
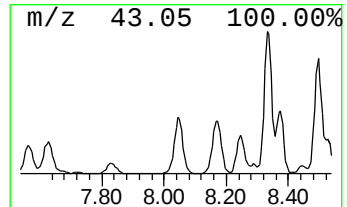
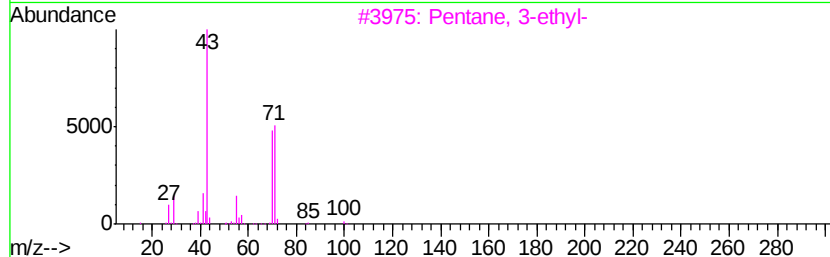
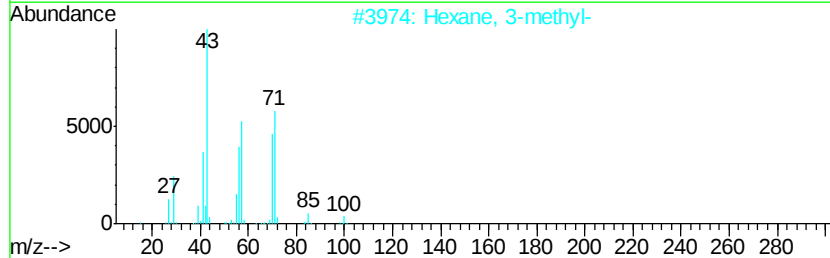
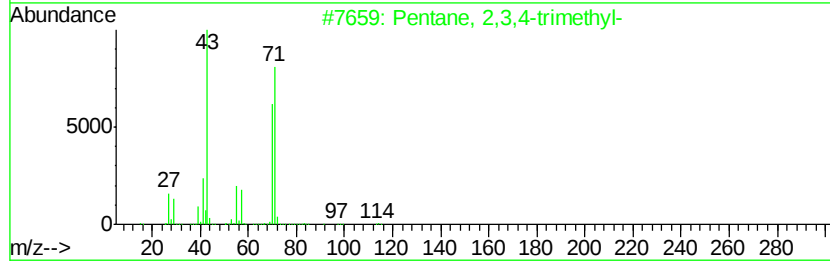
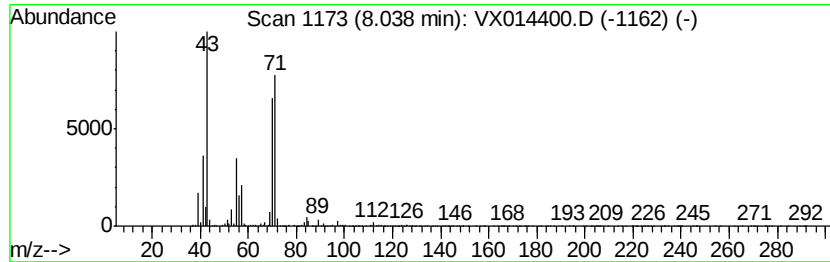
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	24.25 ug/l	893242	1,4-Difluorobenzene	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	86
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	78
3		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
4		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	78
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

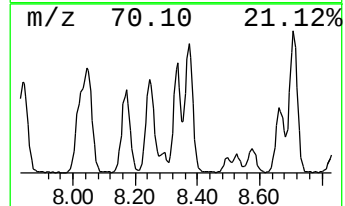
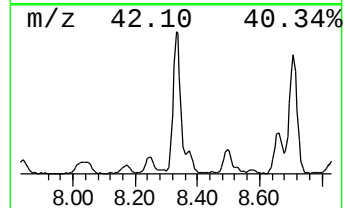
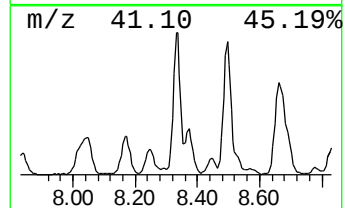
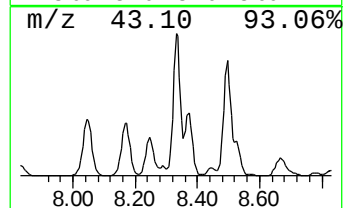
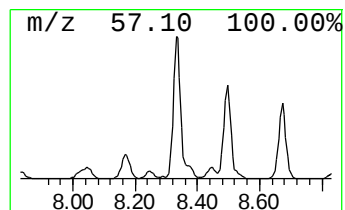
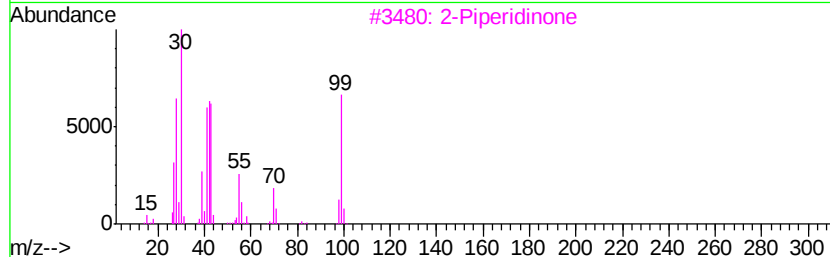
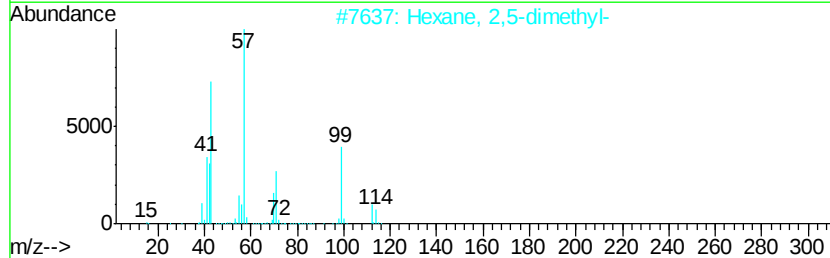
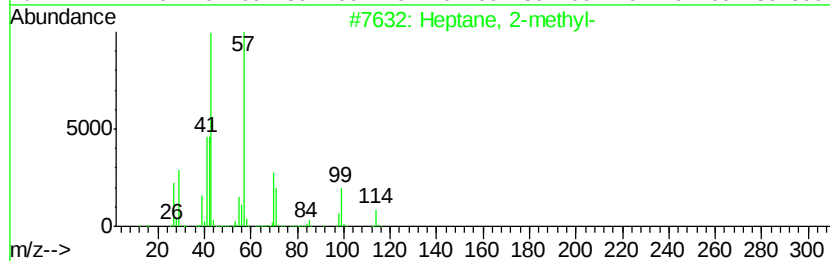
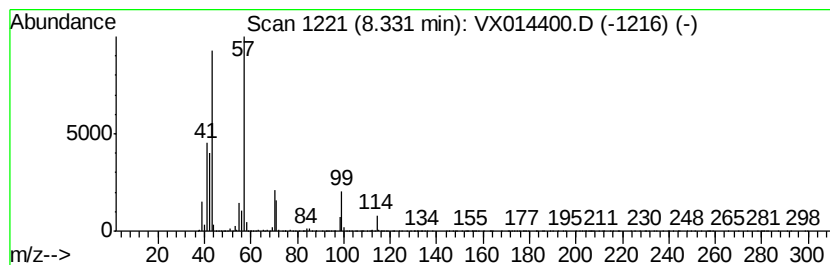
Instrument :
MSVOA_X
ClientSampleId :
WP022SB454_191227-12-14

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 6 Heptane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.33	38.76 ug/l	1428030	1,4-Difluorobenzene	6.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-methyl-	114	C8H18	000592-27-8	95
2	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	68
3	2-Piperidinone	99	C5H9NO	000675-20-7	58
4	Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	38
5	2,5-Hexanedione	114	C6H10O2	000110-13-4	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

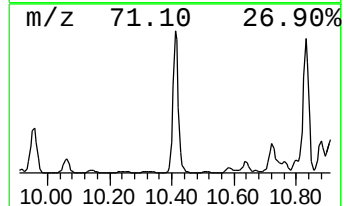
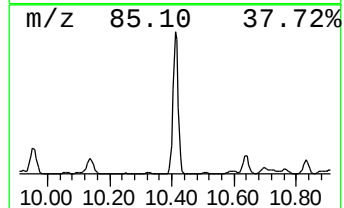
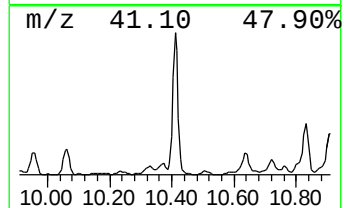
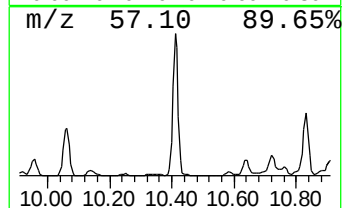
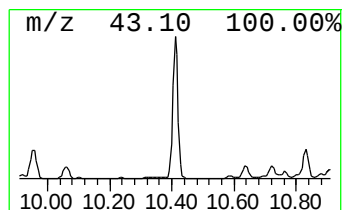
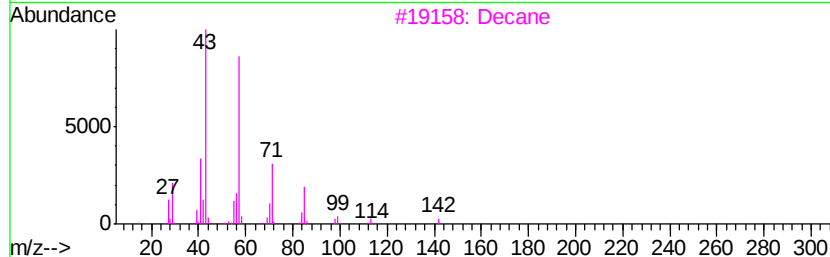
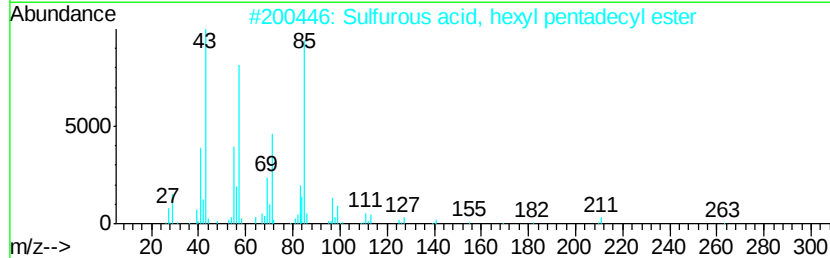
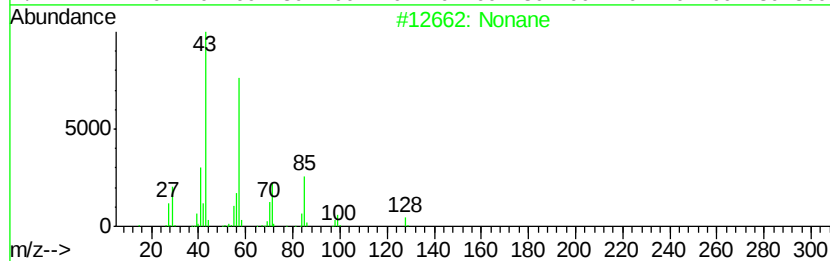
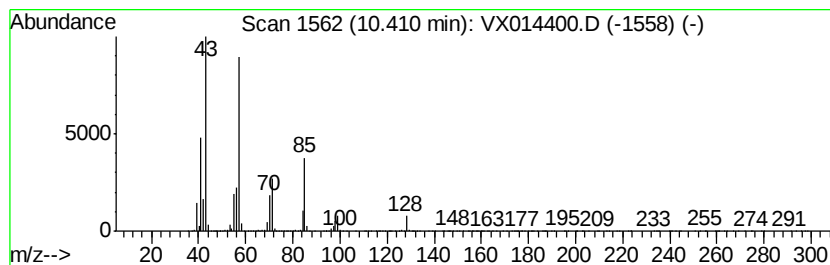
Instrument :
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ClientSampleId :
WP022SB454_191227-12-14

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 7 Nonane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	22.41 ug/l	15454300	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane		128 C9H20	000111-84-2 97
2	Sulfurous acid, hexyl pentadecyl...		376 C21H44O3S	1000309-13-7 72
3	Decane		142 C10H22	000124-18-5 64
4	4,4-Dimethyl octane		142 C10H22	015869-95-1 59
5	Octane, 3,5-dimethyl-		142 C10H22	015869-93-9 53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

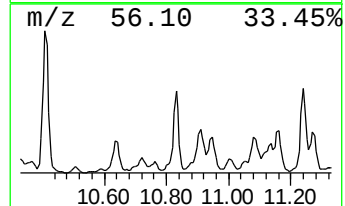
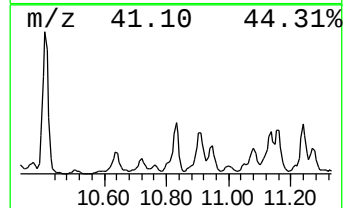
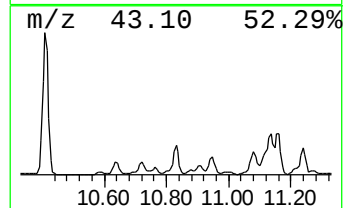
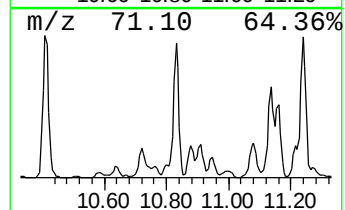
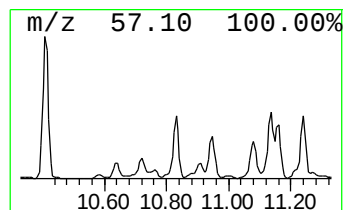
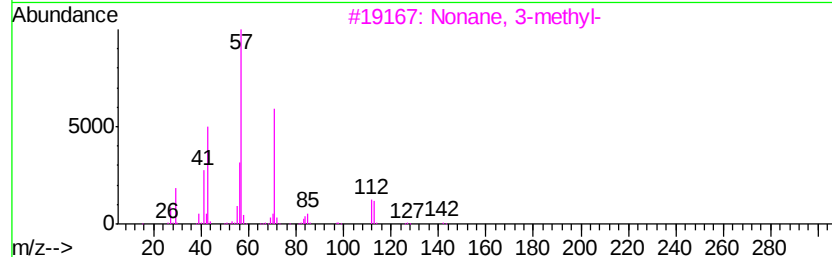
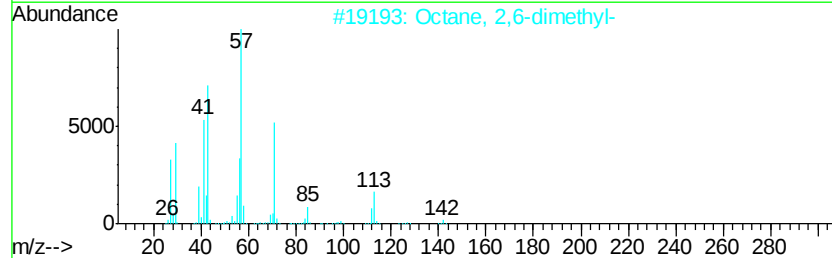
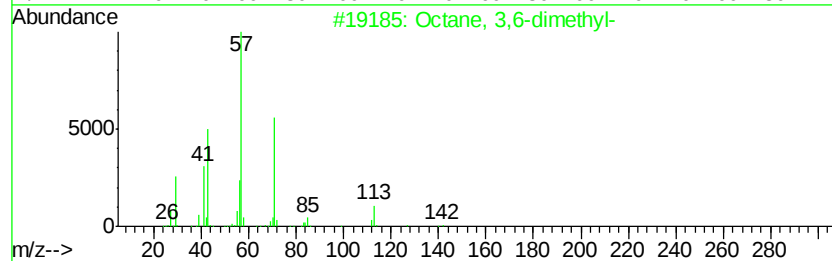
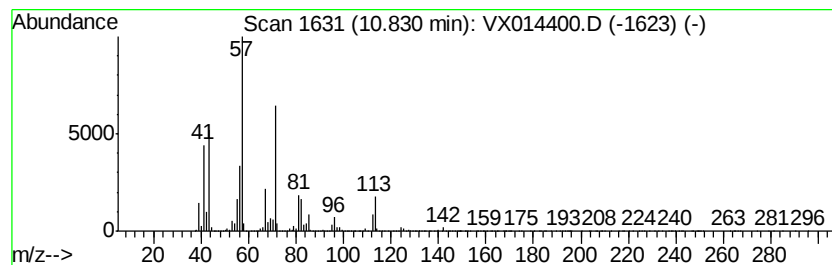
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ClientSampleId :
WP022SB454_191227-12-14

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 8 Octane, 3,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	10.71 ug/l	7384680	Chlorobenzene-d5	10.11
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octane, 3,6-dimethyl-	142 C10H22	015869-94-0	81
2	Octane, 2,6-dimethyl-	142 C10H22	002051-30-1	81
3	Nonane, 3-methyl-	142 C10H22	005911-04-6	76
4	Sulfurous acid, 2-ethylhexyl hex...	278 C14H30O3S	1000309-20-2	52
5	Butane, 2,2-dimethyl-	86 C6H14	000075-83-2	47



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96µ/5mL/100µL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

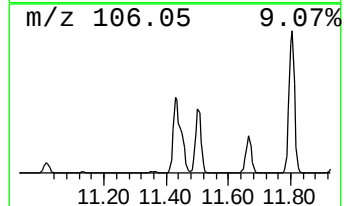
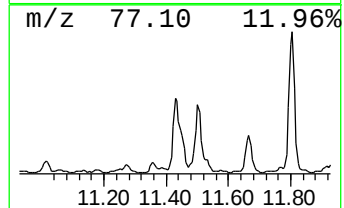
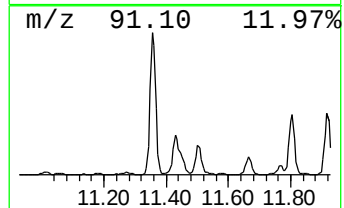
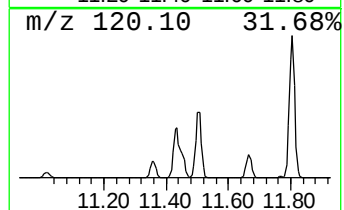
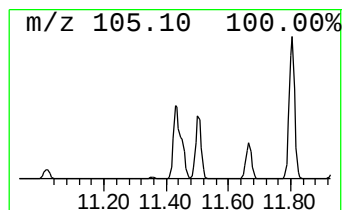
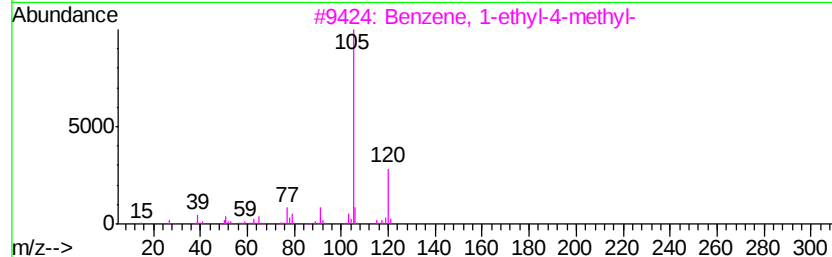
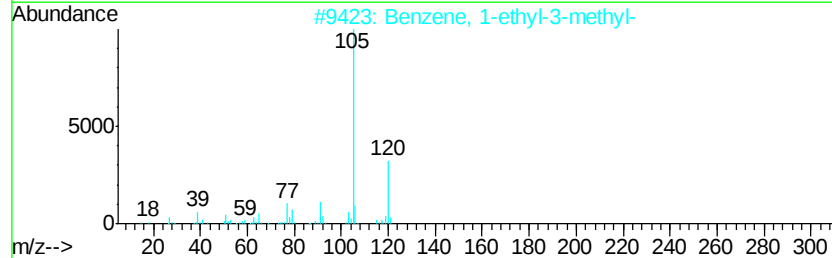
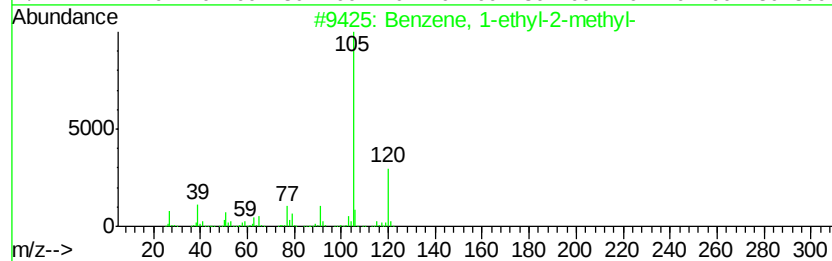
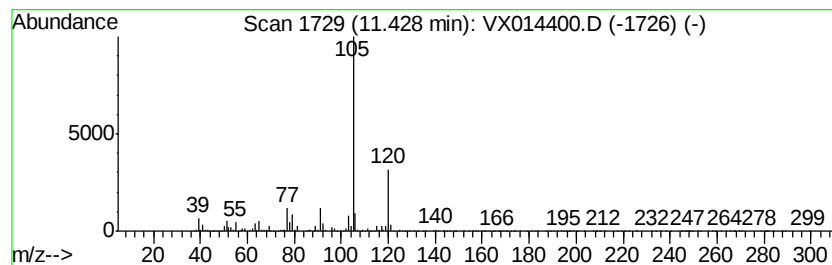
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ClientSampleId :
WP022SB454_191227-12-14

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	11.16 ug/l	6512610	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 94
2		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 94
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 93
4		Mesitylene	120 C9H12	000108-67-8 87
5		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX010320\
 Data File : VX014400.D
 Acq On : 03 Jan 2020 15:04
 Operator : JC/SP
 Sample : K6476-01
 Misc : 6.96µ/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 11 Sample Multiplier: 1

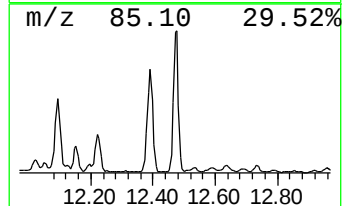
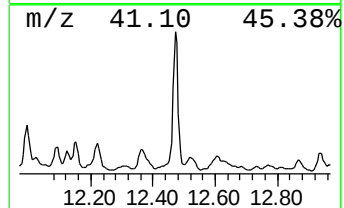
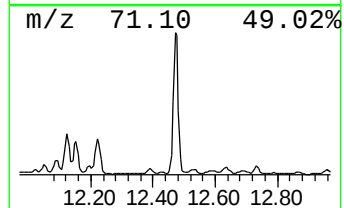
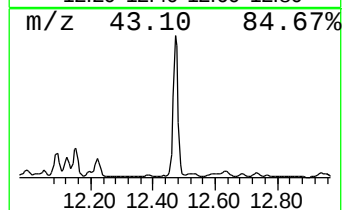
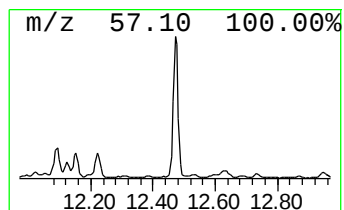
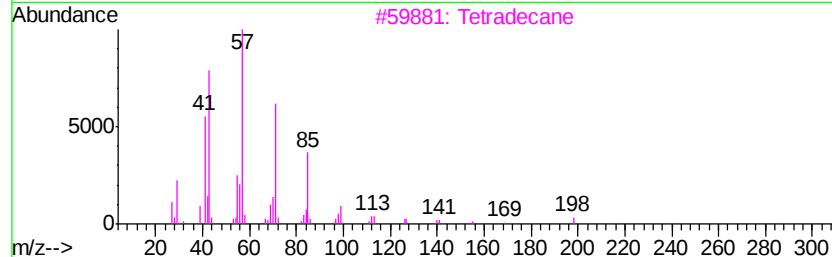
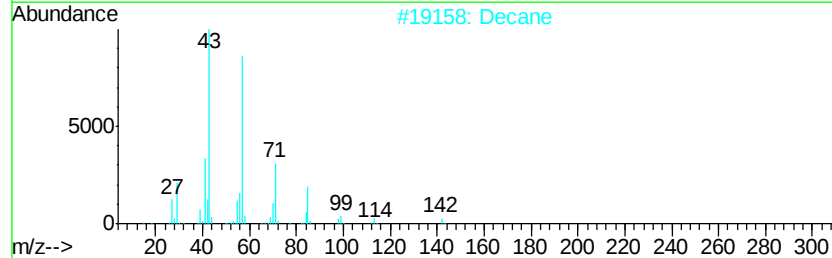
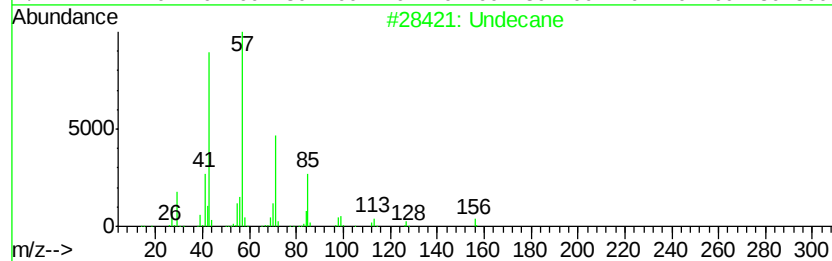
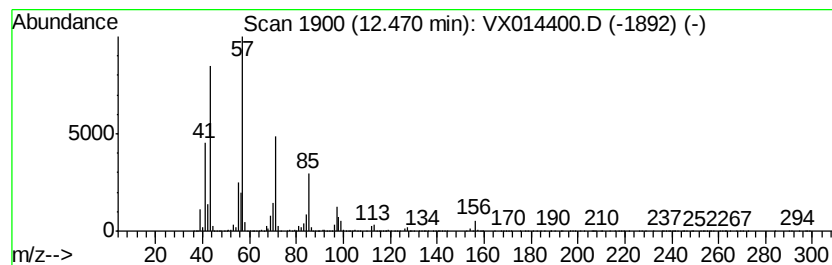
Instrument :
 MSVOA_X
 ClientSampleId :
 WP022SB454_191227-12-14

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.M
 Quant Title : SW846 8260

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

 Peak Number 10 Undecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	24.67 ug/l	14400700	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane	156 C11H24	001120-21-4	94
2	Decane	142 C10H22	000124-18-5	86
3	Tetradecane	198 C14H30	000629-59-4	80
4	Tetracontane, 3,5,24-trimethyl-	605 C43H88	055162-61-3	78
5	Decane, 3,6-dimethyl-	170 C12H26	017312-53-7	74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX010320\
Data File : VX014400.D
Acq On : 03 Jan 2020 15:04
Operator : JC/SP
Sample : K6476-01
Misc : 6.96g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WP022SB454_191227-12-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X121319W.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
n-Hexane	3.50	45.0	ug/l	711583	1	5.65	791315	50.0
Hexane, 3-methyl-	5.91	100.7	ug/l	1593480	1	5.65	791315	50.0
Butane, 2,2,3,3-t...	6.33	42.2	ug/l	1554260	2	6.85	1842080	50.0
Heptane	6.69	72.4	ug/l	2668900	2	6.85	1842080	50.0
Pentane, 2,3,4-tr...	8.04	24.3	ug/l	893242	2	6.85	1842080	50.0
Heptane, 2-methyl-	8.33	38.8	ug/l	1428030	2	6.85	1842080	50.0
Nonane	10.41	22.4	ug/l	15454300	3	10.11	34488400	50.0
Octane, 3,6-dimet...	10.83	10.7	ug/l	7384680	3	10.11	34488400	50.0
Benzene, 1-ethyl-...	11.43	11.2	ug/l	6512610	4	12.08	29184800	50.0
Undecane	12.47	24.7	ug/l	14400700	4	12.08	29184800	50.0