

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091619\
 Data File : VX012416.D
 Acq On : 16 Sep 2019 19:29
 Operator : JC/SP
 Sample : K4830-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S13-0272

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\82X091619W.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	1.070	26	30	37	rBV	582967	691553	4.11%	0.507%
2	1.777	141	146	157	rVB	218039	302004	1.79%	0.221%
3	2.381	235	245	251	rBV	102254	205118	1.22%	0.150%
4	2.832	311	319	323	rBV	215322	483235	2.87%	0.354%
5	2.881	323	327	330	rVV	426555	843769	5.01%	0.618%
6	2.923	330	334	345	rVB	606323	1202663	7.14%	0.881%
7	3.161	362	373	388	rBV	851677	1932453	11.48%	1.416%
8	4.301	549	560	565	rBV	67220	198324	1.18%	0.145%
9	4.393	565	575	591	rVB	1867806	5504615	32.69%	4.034%
10	5.228	699	712	733	rVB2	62861	239848	1.42%	0.176%
11	5.490	744	755	758	rBV	102456	260667	1.55%	0.191%
12	5.575	758	769	779	rVV	1759045	5370454	31.89%	3.935%
13	5.716	786	792	806	rVB	257808	794997	4.72%	0.583%
14	5.929	814	827	839	rBV4	380648	1294603	7.69%	0.949%
15	6.057	839	848	855	rBV	113540	290243	1.72%	0.213%
16	6.252	869	880	890	rBV2	339301	1145608	6.80%	0.839%
17	6.350	890	896	904	rVV2	328142	898218	5.33%	0.658%
18	6.447	904	912	929	rVB	551053	1508584	8.96%	1.105%
19	6.850	966	978	987	rBV	190455	471628	2.80%	0.346%
20	7.459	1064	1078	1089	rBV	3950555	9011561	53.52%	6.603%
21	7.630	1101	1106	1115	rVB	102315	205614	1.22%	0.151%
22	7.727	1115	1122	1132	rVB	382633	736074	4.37%	0.539%
23	7.843	1132	1141	1149	rVB2	187083	391085	2.32%	0.287%
24	8.026	1164	1171	1186	rVB2	182498	426497	2.53%	0.313%
25	8.337	1218	1222	1225	rBV	147276	224028	1.33%	0.164%
26	8.380	1225	1229	1235	rVB2	157726	294837	1.75%	0.216%
27	8.496	1244	1248	1251	rBV	131998	205233	1.22%	0.150%
28	8.660	1268	1275	1279	rBV2	764901	1457172	8.65%	1.068%
29	8.709	1279	1283	1297	rVV2	659083	1254364	7.45%	0.919%
30	8.837	1297	1304	1308	rVV	314904	554092	3.29%	0.406%
31	8.904	1312	1315	1322	rVB	215327	351145	2.09%	0.257%
32	9.038	1331	1337	1348	rVB	609298	1047368	6.22%	0.767%
33	9.160	1348	1357	1363	rBV	313888	510562	3.03%	0.374%
34	9.441	1395	1403	1413	rBV	276692	437204	2.60%	0.320%

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35	9.563	1417	1423	1428	rBV4	303020	688927	4.09%	0.505%
36	9.617	1428	1432	1437	rVB	830239	1169343	6.94%	0.857%
37	9.672	1437	1441	1446	rBV	595352	906115	5.38%	0.664%
38	9.873	1469	1474	1479	rBV4	230713	531108	3.15%	0.389%
39	9.953	1484	1487	1491	rVB	140513	179333	1.06%	0.131%
40	10.062	1499	1505	1509	rBV	291075	437579	2.60%	0.321%
41	10.111	1509	1513	1517	rVV	393435	530487	3.15%	0.389%
42	10.184	1517	1525	1529	rVV	288710	479132	2.85%	0.351%
43	10.251	1529	1536	1542	rVV	4436026	6147434	36.51%	4.505%
44	10.331	1544	1549	1551	rVV2	293396	525240	3.12%	0.385%
45	10.355	1551	1553	1559	rVV2	339140	673721	4.00%	0.494%
46	10.404	1559	1561	1573	rVV2	129362	268698	1.60%	0.197%
47	10.642	1592	1600	1607	rBV4	257887	551123	3.27%	0.404%
48	10.831	1623	1631	1636	rBV3	521903	806264	4.79%	0.591%
49	10.910	1636	1644	1648	rBV3	389409	652781	3.88%	0.478%
50	10.946	1648	1650	1655	rVB3	149022	193552	1.15%	0.142%
51	11.013	1655	1661	1666	rBV	1861825	2354441	13.98%	1.725%
52	11.135	1675	1681	1685	rVB	600130	775276	4.60%	0.568%
53	11.276	1696	1704	1707	rVB5	193037	340425	2.02%	0.249%
54	11.355	1712	1717	1722	rBV	2257077	2777390	16.49%	2.035%
55	11.446	1728	1732	1735	rBV2	138096	216399	1.29%	0.159%
56	11.501	1735	1741	1751	rVB	1954357	2717715	16.14%	1.991%
57	11.666	1763	1768	1776	rBV	875328	1258481	7.47%	0.922%
58	11.763	1781	1784	1786	rVV2	273350	305761	1.82%	0.224%
59	11.806	1786	1791	1796	rVB	13583204	16839280	100.00%	12.339%
60	11.940	1810	1813	1818	rVB	484514	625811	3.72%	0.459%
61	12.074	1829	1835	1840	rVB4	568829	1028001	6.10%	0.753%
62	12.141	1840	1846	1851	rVB	7775979	9438386	56.05%	6.916%
63	12.227	1855	1860	1866	rVB2	211913	334725	1.99%	0.245%
64	12.294	1866	1871	1877	rBV	2588478	3226426	19.16%	2.364%
65	12.367	1878	1883	1890	rVB3	1084820	1788325	10.62%	1.310%
66	12.458	1893	1898	1903	rBV2	339469	494338	2.94%	0.362%
67	12.513	1903	1907	1913	rVB	997651	1256316	7.46%	0.921%
68	12.580	1913	1918	1920	rBV	1036551	1349839	8.02%	0.989%
69	12.605	1920	1922	1928	rVV	1137171	1330705	7.90%	0.975%
70	12.666	1928	1932	1935	rVV	1433290	1612311	9.57%	1.181%
71	12.696	1935	1937	1941	rVB	357432	372264	2.21%	0.273%

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72	12.751	1941	1946	1954	rVB3	972336	1801595	10.70%	1.320%
73	12.897	1964	1970	1974	rVB2	829897	1163929	6.91%	0.853%
74	12.970	1974	1982	1986	rBV	948503	1451155	8.62%	1.063%
75	13.013	1986	1989	1996	rVB	1613258	1985331	11.79%	1.455%
76	13.080	1996	2000	2005	rBV4	208126	361485	2.15%	0.265%
77	13.190	2015	2018	2020	rBV2	191477	215720	1.28%	0.158%
78	13.220	2020	2023	2027	rVB	449227	476610	2.83%	0.349%
79	13.263	2027	2030	2033	rVB2	171154	191648	1.14%	0.140%
80	13.306	2033	2037	2039	rBV2	235993	336592	2.00%	0.247%
81	13.342	2039	2043	2048	rVB2	2760837	3764264	22.35%	2.758%
82	13.434	2054	2058	2061	rBV2	184059	256656	1.52%	0.188%
83	13.476	2061	2065	2073	rVB	1598427	2108977	12.52%	1.545%
84	13.556	2073	2078	2081	rBV2	172307	260299	1.55%	0.191%
85	13.635	2087	2091	2096	rBV4	353954	621980	3.69%	0.456%
86	13.720	2102	2105	2110	rVB2	291010	436562	2.59%	0.320%
87	13.830	2114	2123	2130	rVB	3628512	4926445	29.26%	3.610%
88	13.897	2130	2134	2140	rVB2	364696	592554	3.52%	0.434%
89	13.982	2140	2148	2152	rBV2	332646	562236	3.34%	0.412%
90	14.129	2167	2172	2175	rBV3	179295	282983	1.68%	0.207%
91	14.281	2191	2197	2203	rVB2	445664	848100	5.04%	0.621%
92	14.428	2217	2221	2225	rVB2	163129	202333	1.20%	0.148%
93	14.537	2234	2239	2243	rBV2	382487	559854	3.32%	0.410%
94	14.690	2256	2264	2268	rBV	2289918	3099665	18.41%	2.271%
95	14.836	2283	2288	2296	rBV	1666214	2113668	12.55%	1.549%
96	15.269	2356	2359	2363	rVB	183945	234955	1.40%	0.172%
97	15.543	2399	2404	2417	rVB	293412	507073	3.01%	0.372%
98	15.683	2421	2427	2430	rVV	344299	567191	3.37%	0.416%
99	15.720	2430	2433	2442	rVB	245415	363890	2.16%	0.267%
100	15.909	2455	2464	2474	rVB	82353	246845	1.47%	0.181%

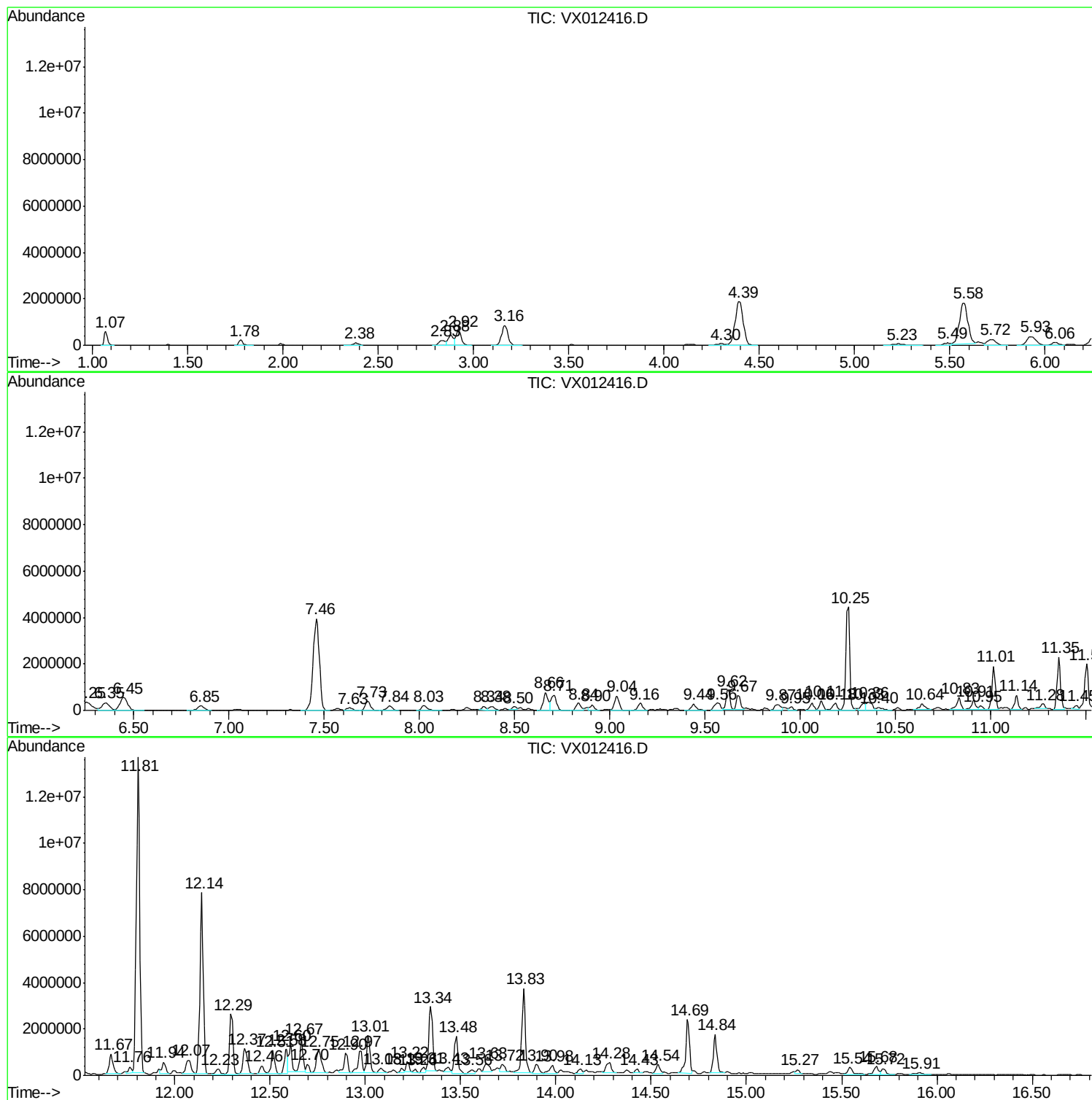
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampled :
S13-0272

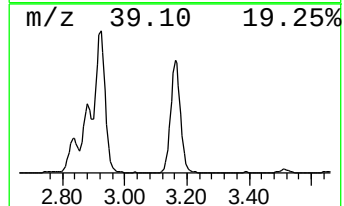
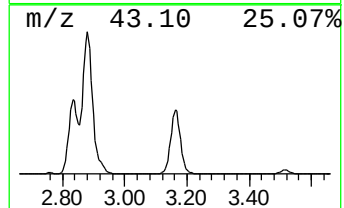
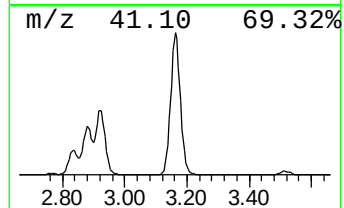
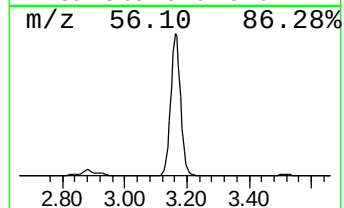
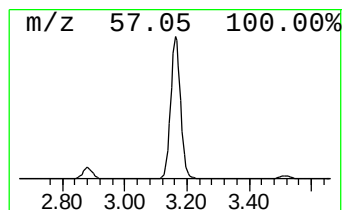
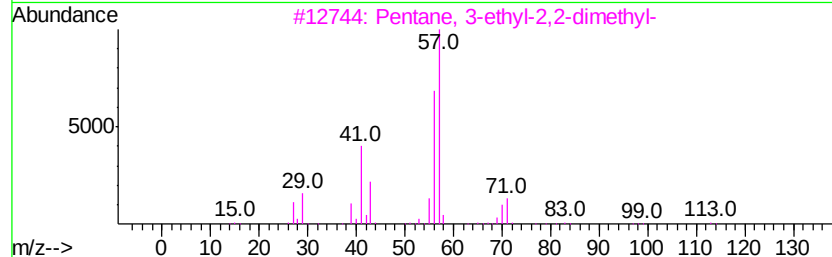
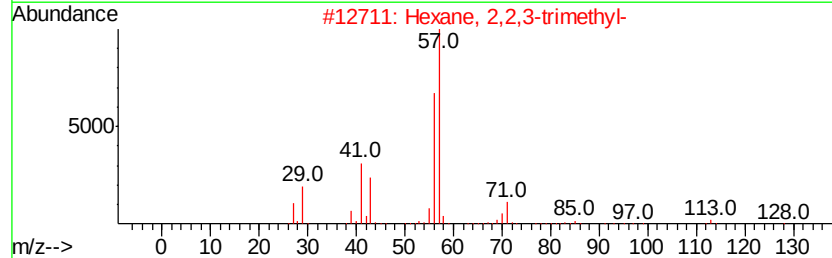
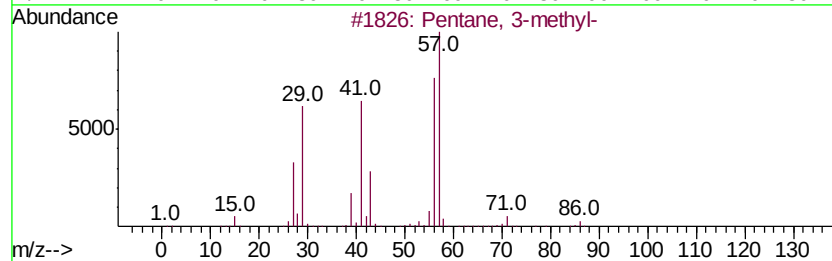
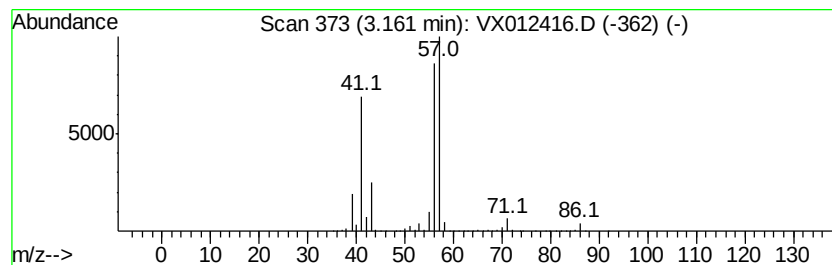
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	121.54 ug/l	1932450	Pentafluorobenzene	5.65

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	43
3		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
4		Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40
5		n-Hexane	86	C6H14	000110-54-3	38



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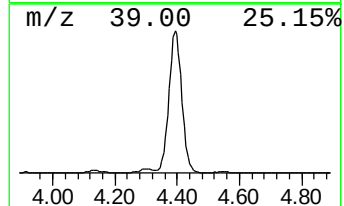
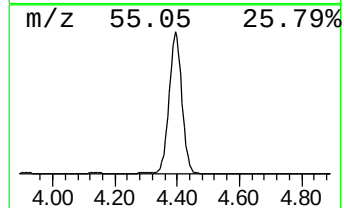
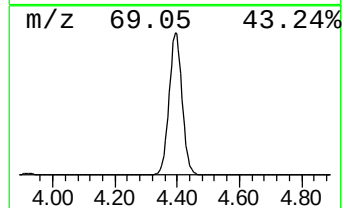
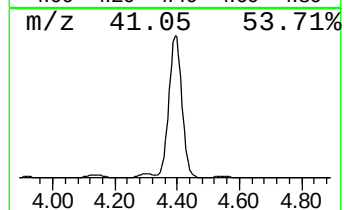
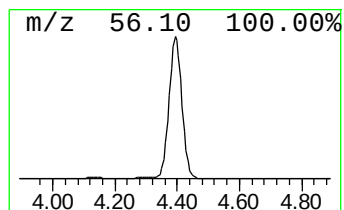
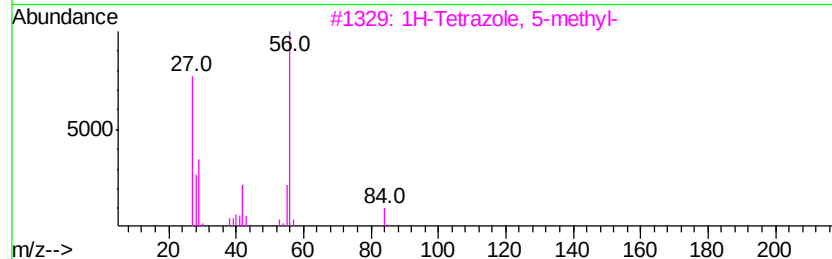
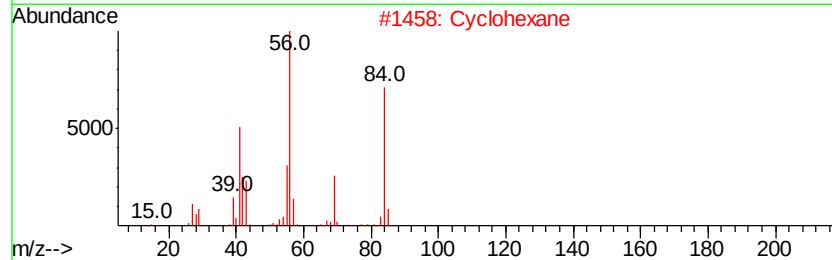
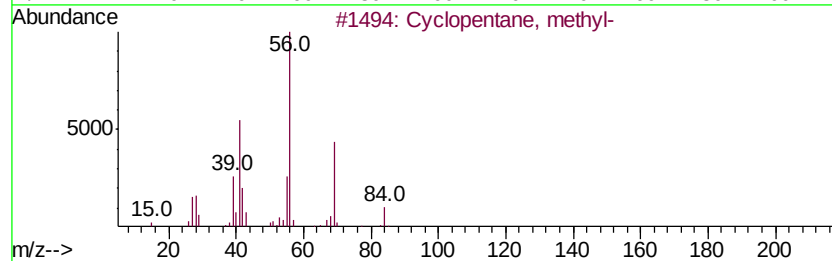
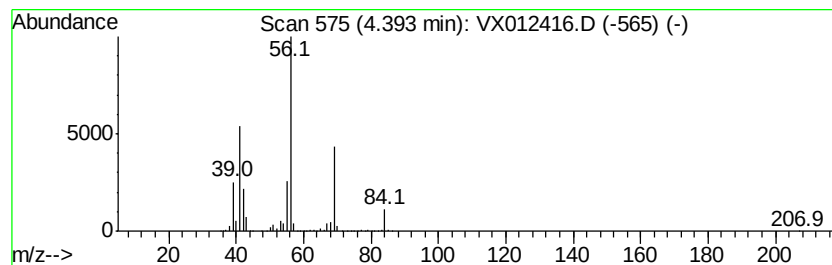
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.39	346.20 ug/l	5504620	Pentafluorobenzene	5.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclobutane	56	C4H8	000287-23-0	53



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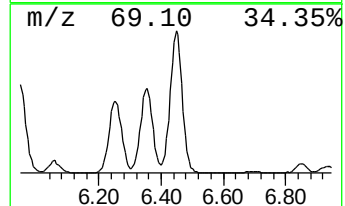
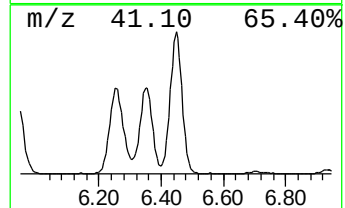
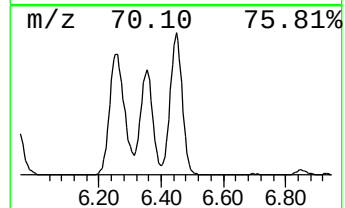
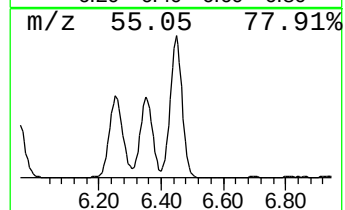
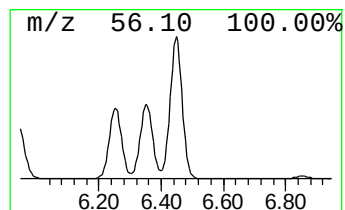
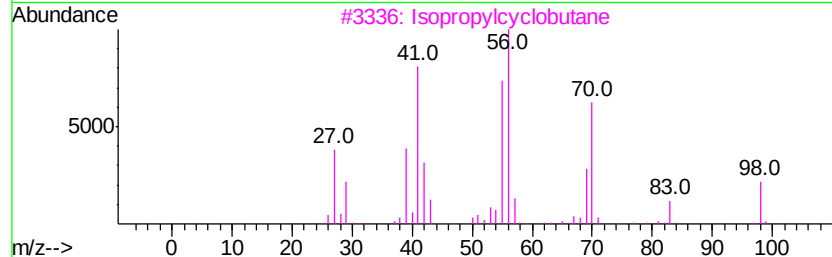
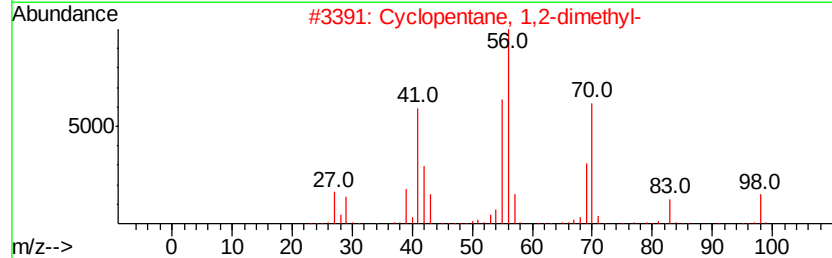
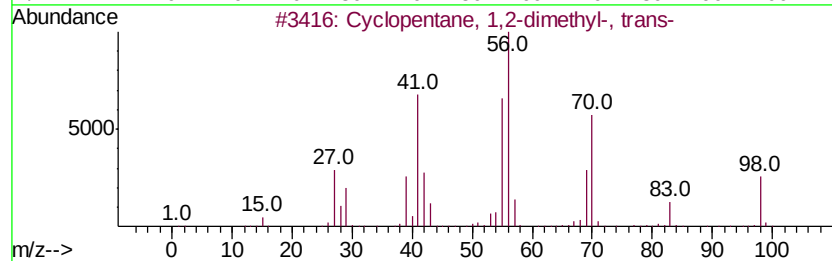
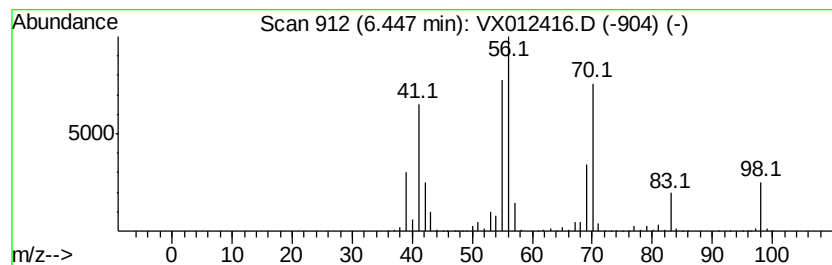
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentane, 1,2-dimethyl-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	159.93 ug/l	1508580	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	94
2		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	94
3		Isopropylcyclobutane	98	C7H14	000872-56-0	94
4		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	86
5		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	86



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0272

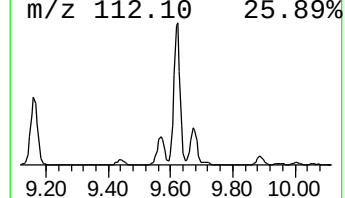
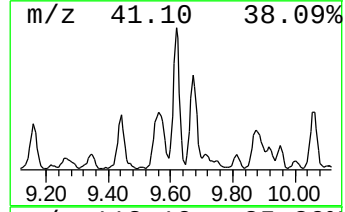
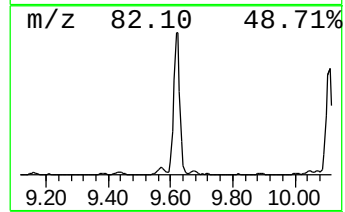
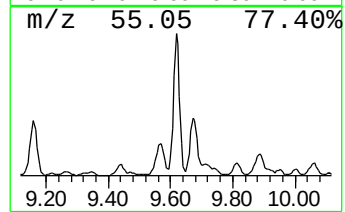
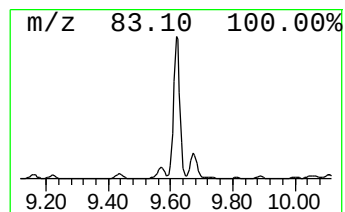
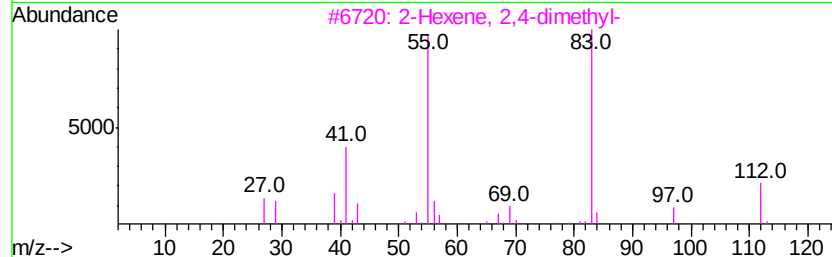
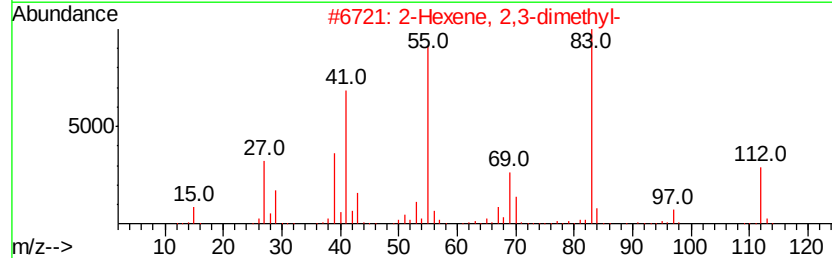
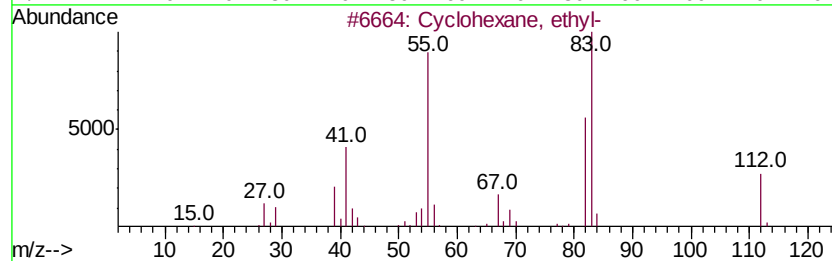
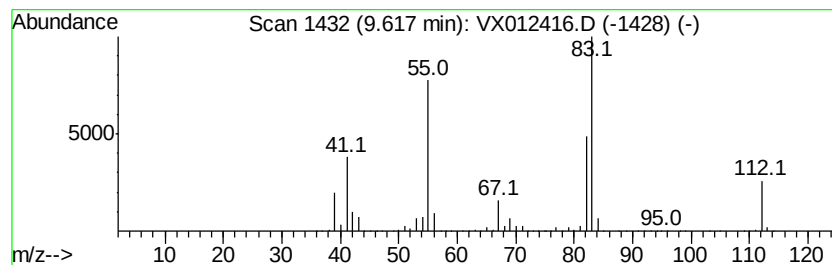
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	110.21 ug/l	1169340	Chlorobenzene-d5	10.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	97
2		2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	64
3		2-Hexene, 2,4-dimethyl-	112	C8H16	014255-23-3	59
4		Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	53
5		2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0272

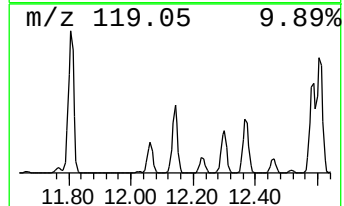
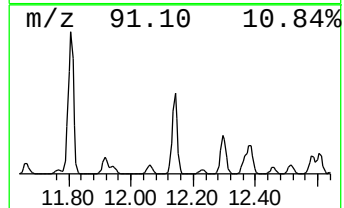
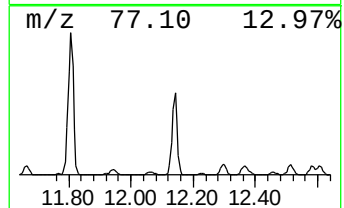
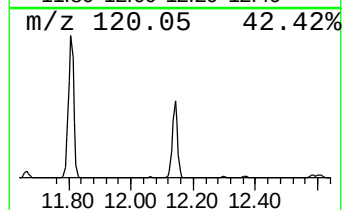
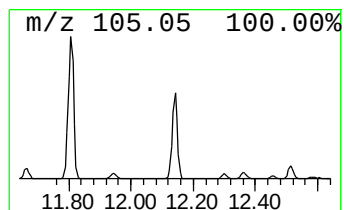
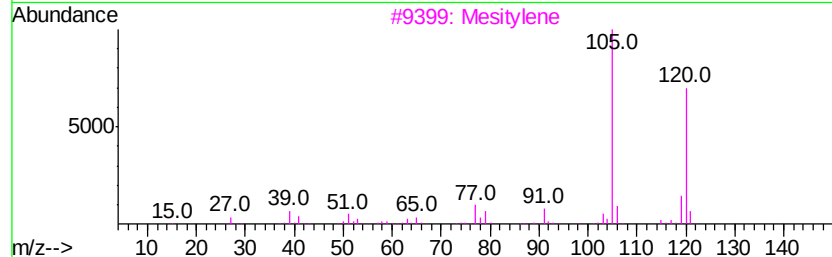
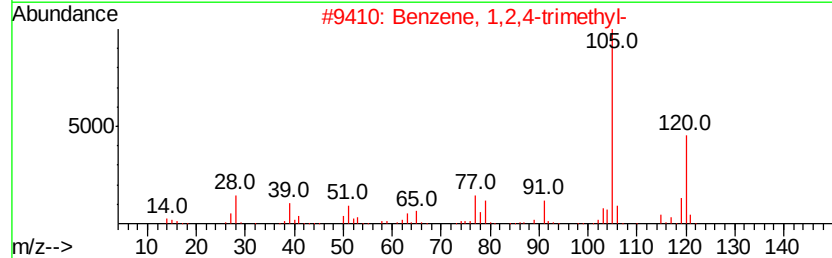
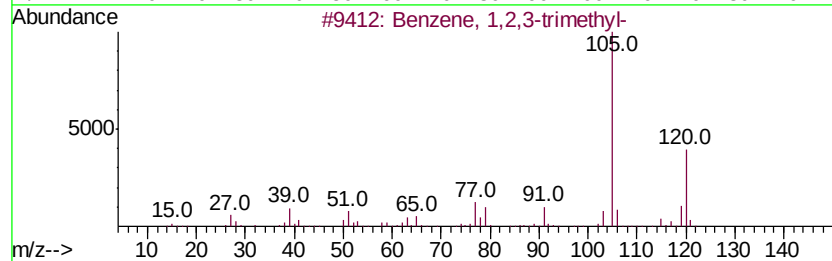
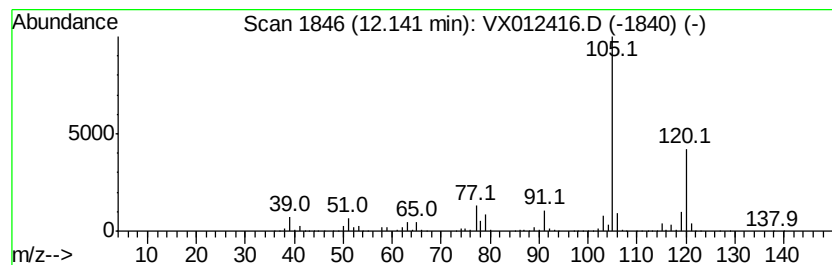
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	459.07 ug/l	9438390	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

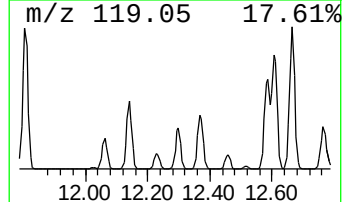
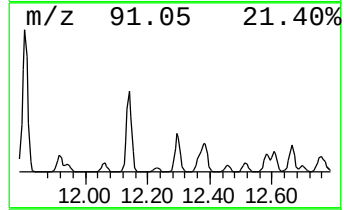
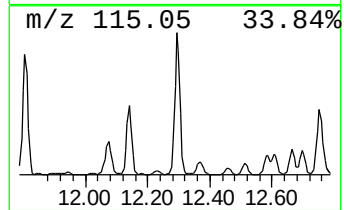
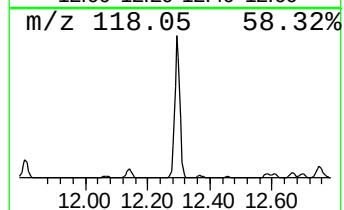
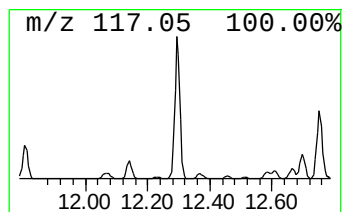
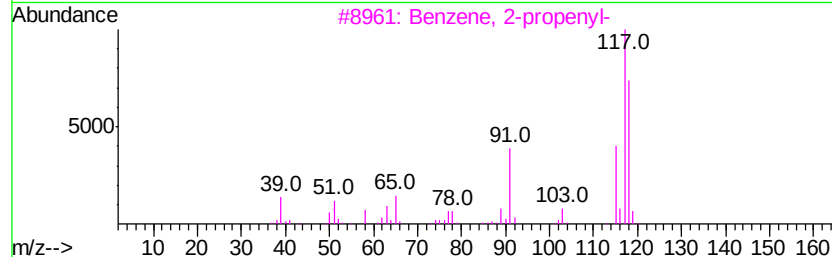
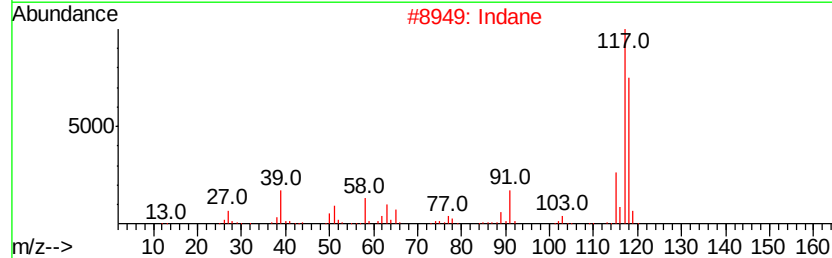
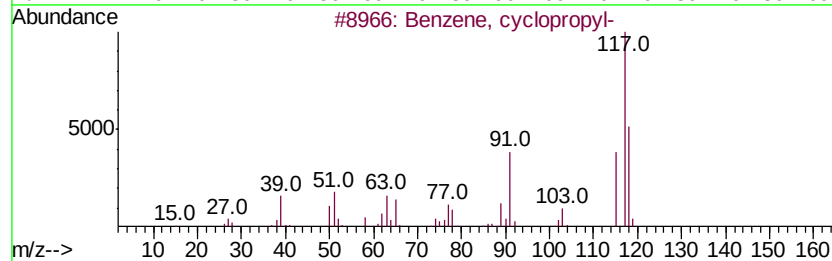
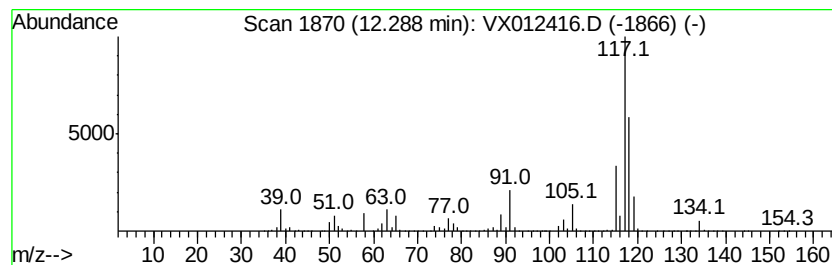
Instrument :
MSVOA_X
ClientSampleId :
S13-0272

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, cyclopropyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.29	156.93 ug/l	3226430	1,4-Dichlorobenzene-d4	12.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cvclopropyl-	118	C9H10	000873-49-4	70
2	Indane	118	C9H10	000496-11-7	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0272

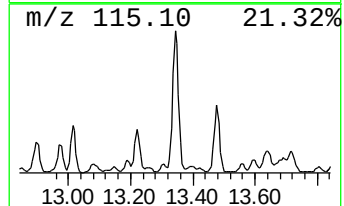
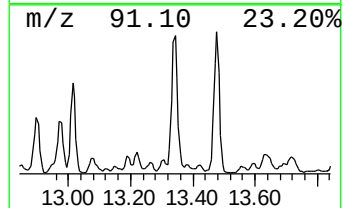
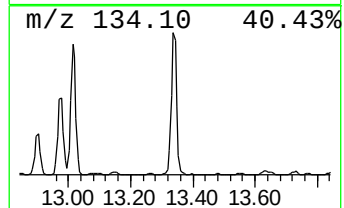
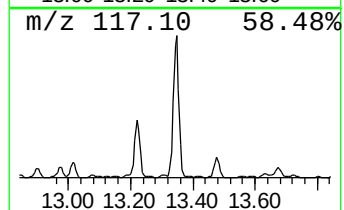
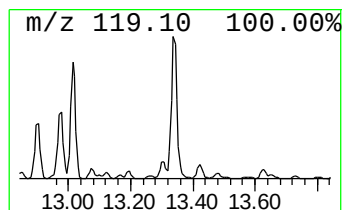
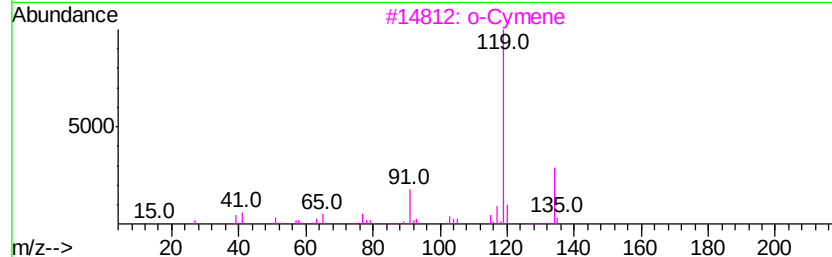
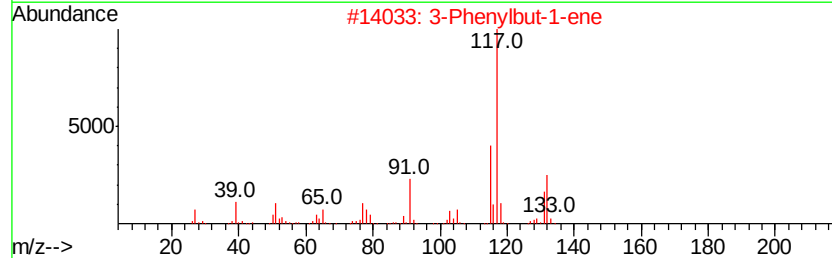
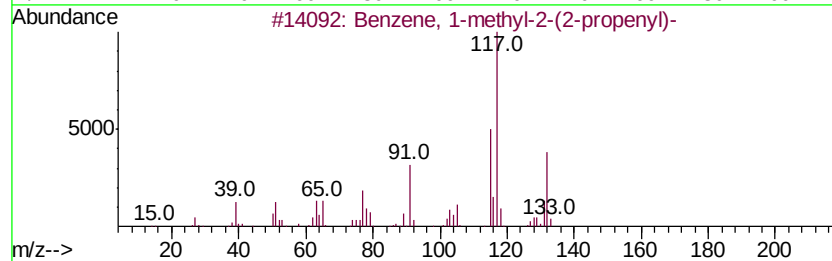
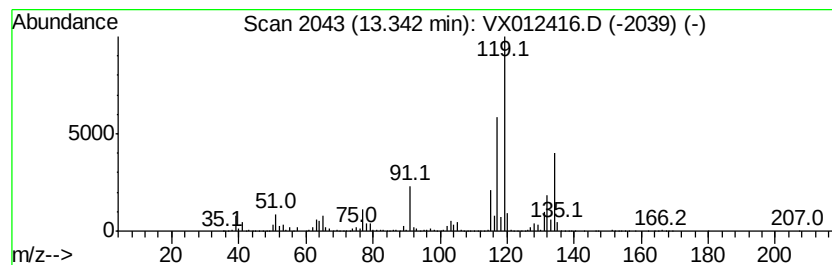
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	183.09 ug/l	3764260	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		o-Cymene	134	C10H14	000527-84-4	64
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
5		p-Cymene	134	C10H14	000099-87-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

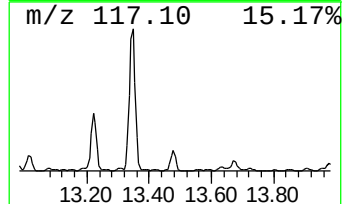
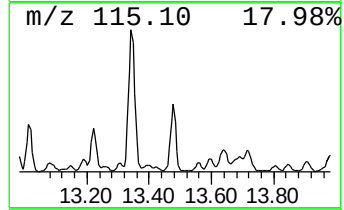
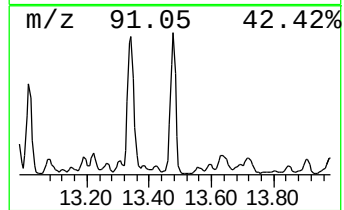
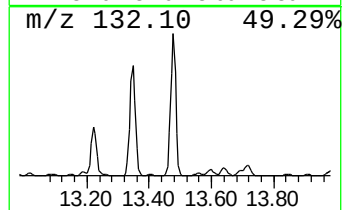
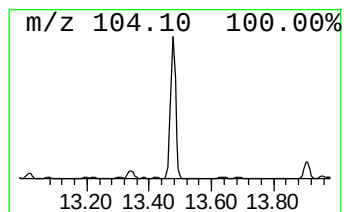
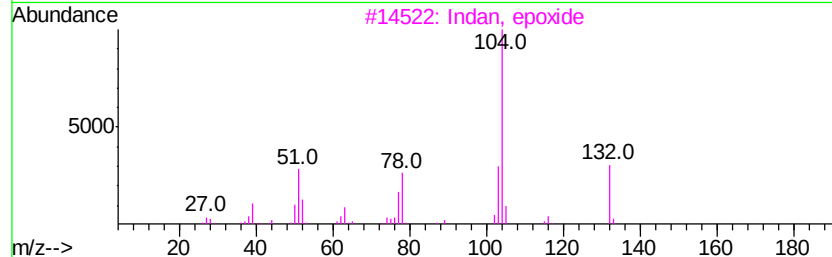
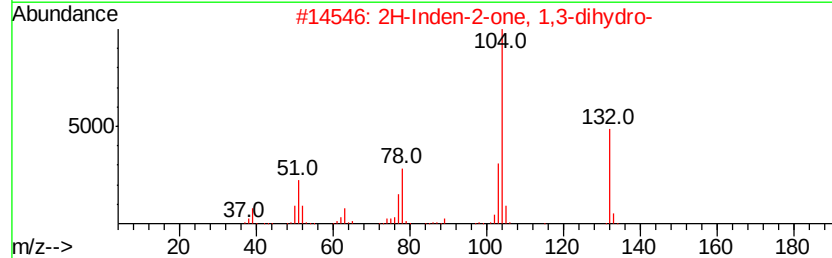
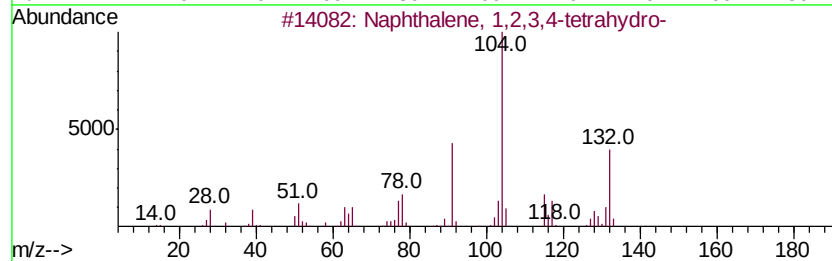
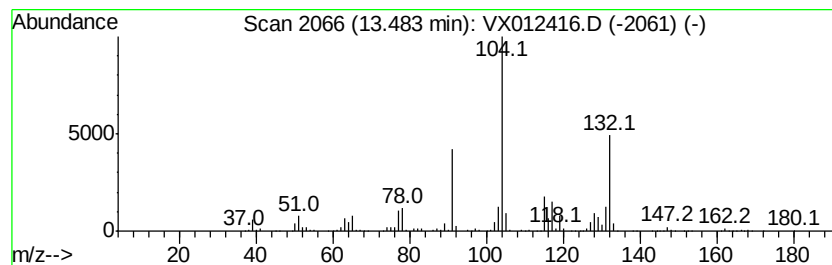
Instrument :
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ClientSampleId :
S13-0272

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.M
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 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	102.58 ug/l	2108980	1,4-Dichlorobenzene-d4	12.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 96
2		2H-Inden-2-one, 1,3-dihydro-	132 C9H8O	000615-13-4 53
3		Indan, epoxide	132 C9H8O	000768-22-9 52
4		Benzene, 3-cyclohexen-1-yl-	158 C12H14	004994-16-5 47
5		Benzene, cyclobutyl-	132 C10H12	004392-30-7 45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
S13-0272

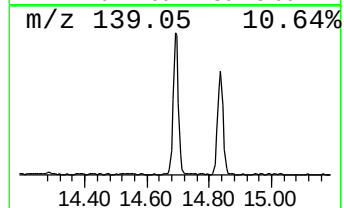
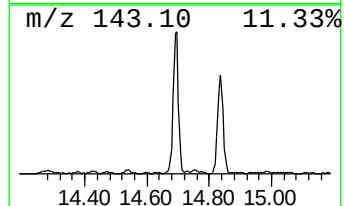
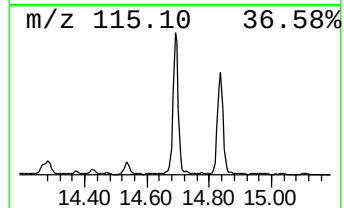
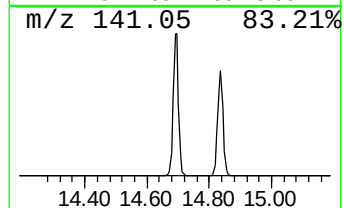
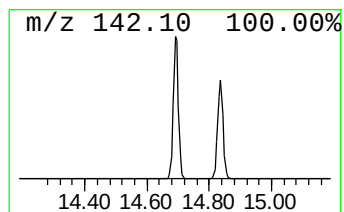
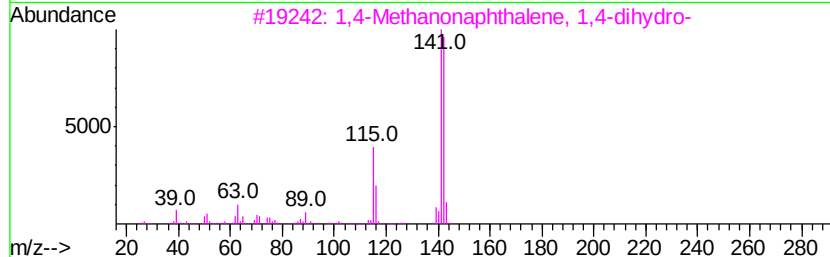
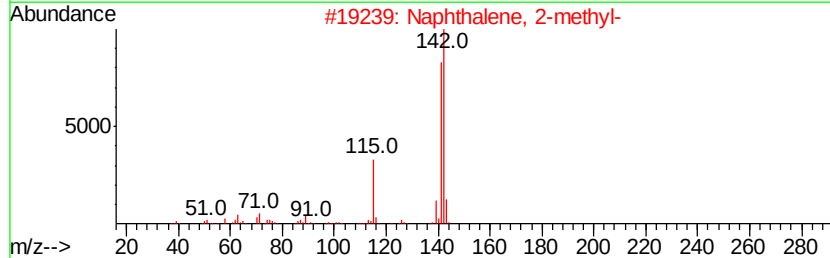
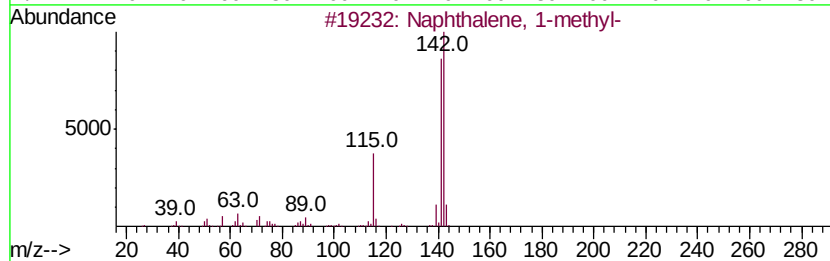
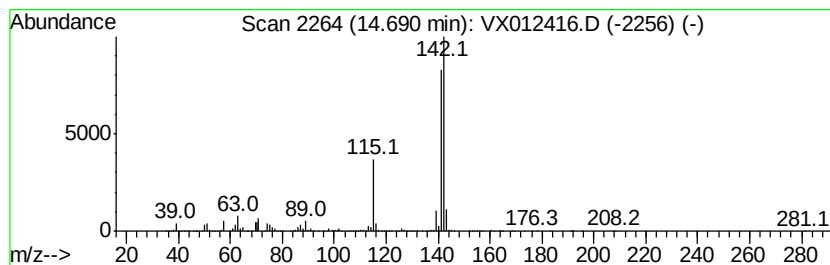
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	150.76 ug/l	3099670	1,4-Dichlorobenzene-d4	12.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	93
4		Benzocycloheptatriene	142	C11H10	000264-09-5	90
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX091619\
Data File : VX012416.D
Acq On : 16 Sep 2019 19:29
Operator : JC/SP
Sample : K4830-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S13-0272

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X091619W.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, 3-methyl-	3.16	121.5	ug/l	1932450	1	5.65	794997	50.0
Cyclopentane, met...	4.39	346.2	ug/l	5504620	1	5.65	794997	50.0
Cyclopentane, 1,2...	6.45	159.9	ug/l	1508580	2	6.86	471628	50.0
Cyclohexane, ethyl-	9.62	110.2	ug/l	1169340	3	10.11	530487	50.0
Benzene, 1,2,3-tr...	12.14	459.1	ug/l	9438390	4	12.07	1028000	50.0
Benzene, cyclopro...	12.29	156.9	ug/l	3226430	4	12.07	1028000	50.0
Benzene, 1-methyl...	13.34	183.1	ug/l	3764260	4	12.07	1028000	50.0
Naphthalene, 1,2,...	13.48	102.6	ug/l	2108980	4	12.07	1028000	50.0
Naphthalene, 1-me...	14.69	150.8	ug/l	3099670	4	12.07	1028000	50.0