

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
 Data File : VX027173.D
 Acq On : 24 Feb 2022 16:26
 Operator : JC/MD
 Sample : N1576-01
 Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PE-4

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
 Title : SW846 8260

Signal : TIC: VX027173.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.800	276	281	293	rVB	110462	227313	2.57%	0.269%
2	3.081	318	327	342	rVB	87524	193654	2.19%	0.229%
3	3.428	375	384	396	rBV	74034	165134	1.87%	0.195%
4	4.196	500	510	518	rBV	59225	173774	1.97%	0.206%
5	4.288	518	525	537	rVB	73672	206962	2.34%	0.245%
6	5.391	689	706	713	rBV3	51646	184426	2.09%	0.218%
7	5.495	713	723	730	rBV	203526	642069	7.26%	0.760%
8	5.611	737	742	763	rVB	165716	536170	6.06%	0.634%
9	5.818	763	776	790	rBV	266630	772181	8.73%	0.914%
10	5.958	790	799	813	rVB	51168	143237	1.62%	0.169%
11	6.153	813	831	837	rBV2	80241	278792	3.15%	0.330%
12	6.245	837	846	856	rVV	412586	1273795	14.41%	1.507%
13	6.355	856	864	874	rVB2	63419	185054	2.09%	0.219%
14	6.605	892	905	913	rBV	216154	531700	6.01%	0.629%
15	6.763	924	931	937	rBV	134000	350485	3.96%	0.415%
16	6.842	937	944	951	rVB3	103572	320030	3.62%	0.379%
17	7.172	988	998	1006	rBV3	63383	157461	1.78%	0.186%
18	7.379	1019	1032	1043	rBV3	263517	715912	8.10%	0.847%
19	7.495	1043	1051	1056	rBV	135146	276813	3.13%	0.327%
20	7.556	1056	1061	1072	rVV	160506	396565	4.48%	0.469%
21	7.653	1072	1077	1089	rVB	89267	179699	2.03%	0.213%
22	7.769	1089	1096	1104	rVB2	79720	164740	1.86%	0.195%
23	7.982	1121	1131	1143	rVB	334464	741770	8.39%	0.878%
24	8.104	1143	1151	1160	rBV2	389981	952604	10.77%	1.127%
25	8.184	1160	1164	1170	rVV2	192762	396934	4.49%	0.470%
26	8.275	1174	1179	1183	rVV	396381	719543	8.14%	0.851%
27	8.312	1183	1185	1191	rVV2	198325	279819	3.16%	0.331%
28	8.434	1200	1205	1214	rVV	425911	834835	9.44%	0.988%
29	8.507	1214	1217	1227	rVB4	85777	187230	2.12%	0.222%
30	8.610	1227	1234	1237	rBV3	248652	554375	6.27%	0.656%
31	8.647	1237	1240	1248	rVB	289498	518589	5.86%	0.614%
32	8.775	1248	1261	1264	rBV5	115459	298574	3.38%	0.353%
33	8.818	1264	1268	1277	rVV3	120906	313868	3.55%	0.371%
34	8.915	1277	1284	1291	rVV3	373406	754860	8.54%	0.893%
35	8.982	1291	1295	1305	rVV4	96304	269024	3.04%	0.318%
36	9.104	1311	1315	1321	rVV2	121032	249399	2.82%	0.295%

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37	9.159	1321	1324	1328	rVV2	113603	184713	2.09%	0.219%
38	9.202	1328	1331	1341	rVB5	65748	144086	1.63%	0.170%
39	9.385	1355	1361	1366	rBV2	122749	201244	2.28%	0.238%
40	9.500	1375	1380	1387	rBV3	224122	449050	5.08%	0.531%

41	9.763	1416	1423	1428	rBV3	78006	154849	1.75%	0.183%
42	9.823	1428	1433	1437	rBV2	86778	180333	2.04%	0.213%
43	9.903	1442	1446	1451	rVB	521034	756530	8.56%	0.895%
44	10.006	1458	1463	1468	rVB	410558	553874	6.26%	0.655%
45	10.061	1468	1472	1476	rBV	267580	371226	4.20%	0.439%

46	10.195	1489	1494	1499	rVB	891303	1222346	13.82%	1.446%
47	10.262	1499	1505	1506	rVV3	86135	139410	1.58%	0.165%
48	10.305	1506	1512	1518	rVV	4041178	5747733	65.00%	6.800%
49	10.360	1518	1521	1533	rVB2	418782	698067	7.89%	0.826%
50	10.586	1553	1558	1562	rBV	166668	224303	2.54%	0.265%

51	10.647	1562	1568	1577	rVV	1171265	1695833	19.18%	2.006%
52	10.781	1582	1590	1594	rBV	159020	267490	3.03%	0.316%
53	10.860	1594	1603	1607	rBV2	135231	247057	2.79%	0.292%
54	10.964	1614	1620	1625	rBV	300200	420885	4.76%	0.498%
55	11.031	1628	1631	1634	rVV	108792	147813	1.67%	0.175%

56	11.085	1634	1640	1642	rVV2	443324	668821	7.56%	0.791%
57	11.110	1642	1644	1651	rVV2	326285	508698	5.75%	0.602%
58	11.195	1651	1658	1665	rVV	265377	415785	4.70%	0.492%
59	11.305	1665	1676	1683	rVV	1005537	1454179	16.45%	1.720%
60	11.384	1683	1689	1696	rVV	4865274	8512486	96.27%	10.071%

61	11.457	1696	1701	1712	rVB	2462447	3522757	39.84%	4.168%
62	11.616	1722	1727	1734	rBV	1799411	2277196	25.75%	2.694%
63	11.756	1745	1750	1758	rVB	6680286	8842556	100.00%	10.461%
64	11.890	1770	1772	1776	rVB	142044	168671	1.91%	0.200%
65	11.976	1782	1786	1789	rVB	278302	319835	3.62%	0.378%

66	12.024	1789	1794	1800	rBV3	338555	577740	6.53%	0.683%
67	12.091	1800	1805	1810	rBV	1702498	2227594	25.19%	2.635%
68	12.177	1816	1819	1823	rVB2	158236	175806	1.99%	0.208%
69	12.250	1826	1831	1833	rBV2	1294981	1921006	21.72%	2.273%
70	12.268	1833	1834	1838	rVV	1332907	1332591	15.07%	1.577%

71	12.323	1838	1843	1849	rVB2	2017276	3103554	35.10%	3.672%
72	12.427	1849	1860	1863	rBV2	308830	565201	6.39%	0.669%
73	12.463	1863	1866	1873	rVB	497976	657160	7.43%	0.777%
74	12.536	1873	1878	1880	rBV	1036404	1458410	16.49%	1.725%
75	12.555	1880	1881	1887	rVV	951442	974704	11.02%	1.153%

76	12.616	1887	1891	1895	rVV	1745694	2075535	23.47%	2.455%
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77	12.646	1895	1896	1900	rVB	208792	174484	1.97%	0.206%
78	12.701	1900	1905	1914	rVB3	543952	988463	11.18%	1.169%
79	12.847	1925	1929	1934	rVB	335464	420075	4.75%	0.497%
80	12.921	1934	1941	1945	rBV	880843	1269137	14.35%	1.501%
81	12.963	1945	1948	1954	rVB	1268550	1586358	17.94%	1.877%
82	13.024	1954	1958	1960	rBV	232661	322186	3.64%	0.381%
83	13.049	1960	1962	1964	rVV2	255643	284225	3.21%	0.336%
84	13.073	1964	1966	1969	rVB	203396	187189	2.12%	0.221%
85	13.170	1975	1982	1986	rVB3	791851	1103245	12.48%	1.305%
86	13.213	1986	1989	1992	rVB2	178542	194840	2.20%	0.231%
87	13.262	1992	1997	1999	rBV3	294106	466903	5.28%	0.552%
88	13.292	1999	2002	2008	rVV2	1297567	1860896	21.04%	2.202%
89	13.372	2012	2015	2020	rVB	214639	255083	2.88%	0.302%
90	13.427	2020	2024	2032	rVB2	166679	268993	3.04%	0.318%
91	13.506	2032	2037	2040	rBV2	162638	232774	2.63%	0.275%
92	13.542	2040	2043	2046	rVV	242887	312263	3.53%	0.369%
93	13.585	2046	2050	2056	rVV3	345975	647897	7.33%	0.766%
94	13.664	2060	2063	2069	rVB2	214658	368635	4.17%	0.436%
95	13.774	2077	2081	2088	rVB	624368	828495	9.37%	0.980%
96	13.926	2100	2106	2111	rBV4	138330	263904	2.98%	0.312%
97	14.079	2127	2131	2135	rBV	179096	227623	2.57%	0.269%
98	14.213	2149	2153	2163	rVV	154122	259807	2.94%	0.307%
99	14.640	2213	2223	2233	rVB	673395	904184	10.23%	1.070%
100	14.780	2240	2246	2255	rVB	299867	384779	4.35%	0.455%

Sum of corrected areas: 84526955

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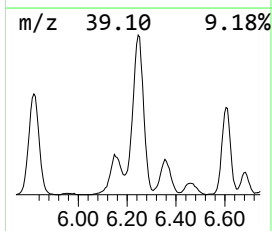
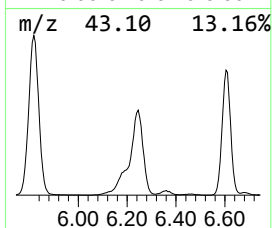
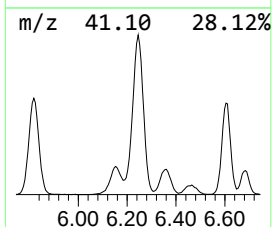
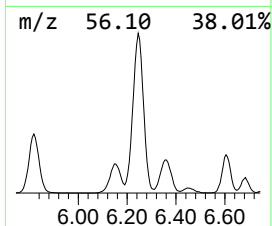
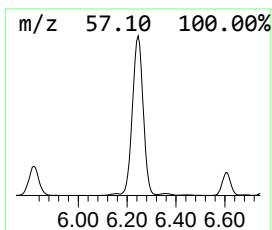
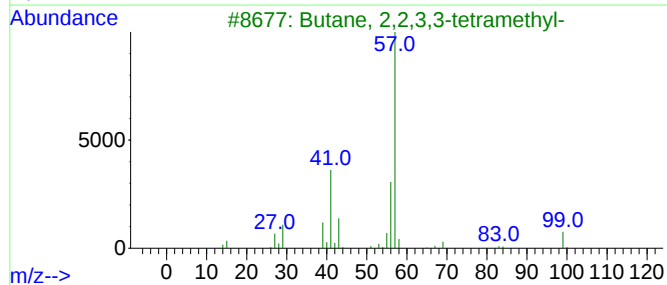
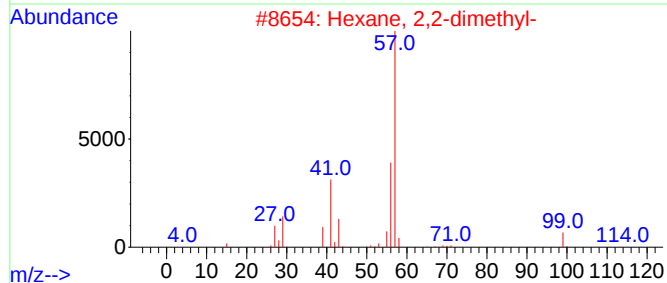
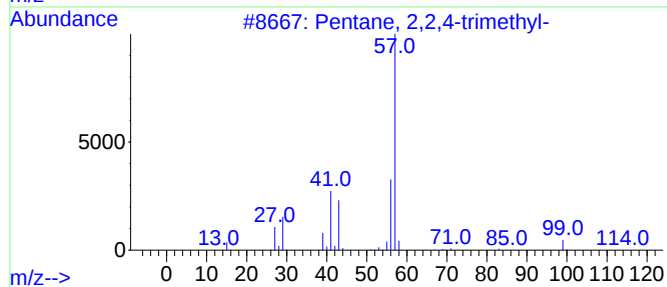
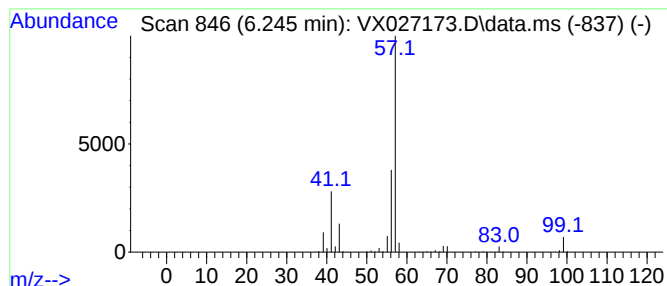
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Pentane, 2,2,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.245	181.72 ug/l	1273800	1,4-Difluorobenzene	6.763

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



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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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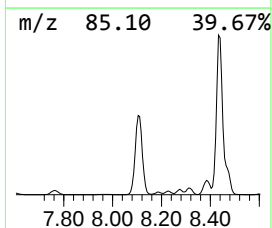
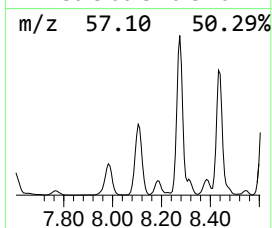
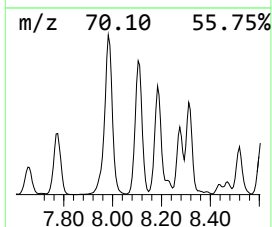
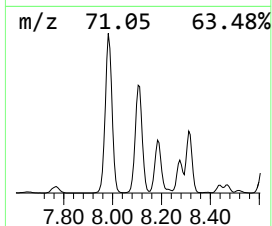
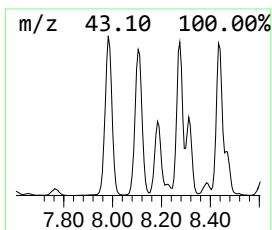
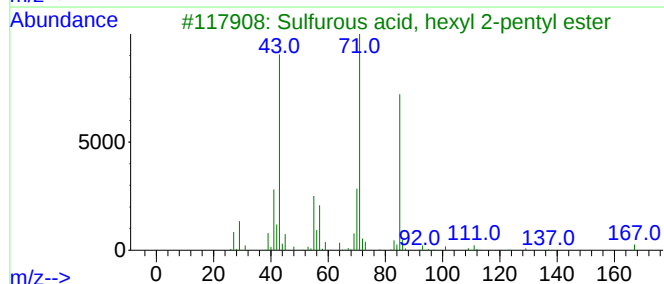
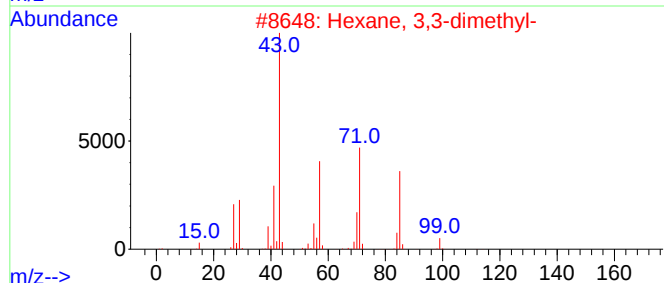
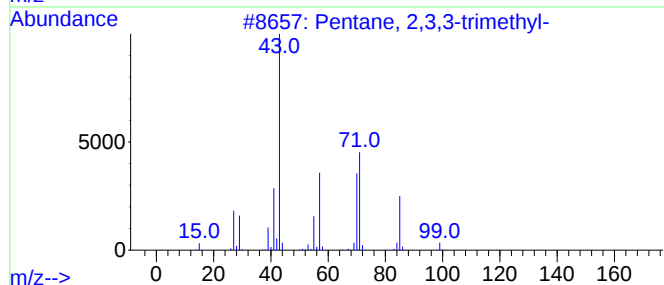
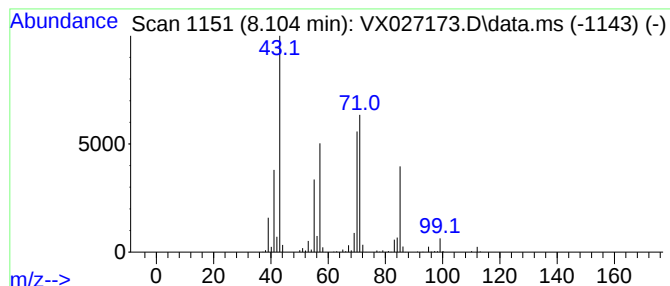
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	135.90 ug/l	952604	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	50
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	50
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	50



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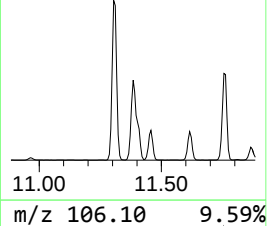
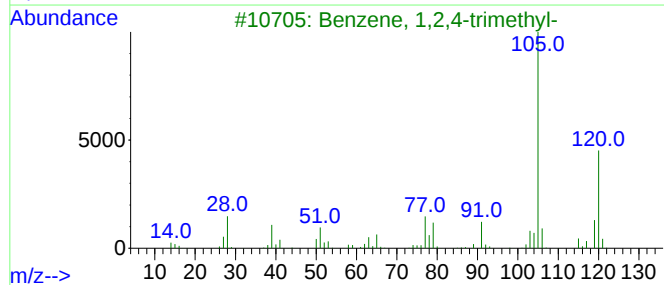
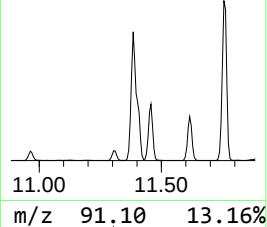
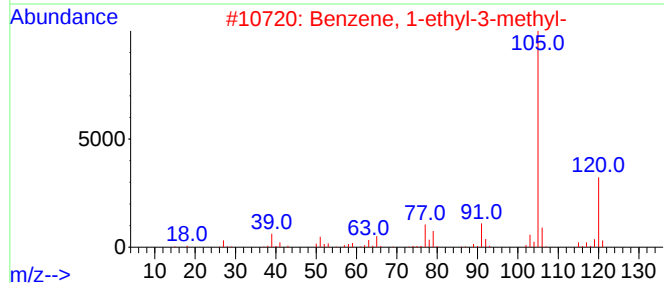
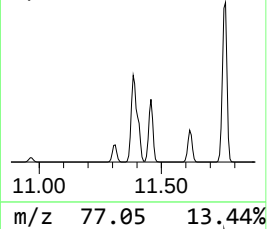
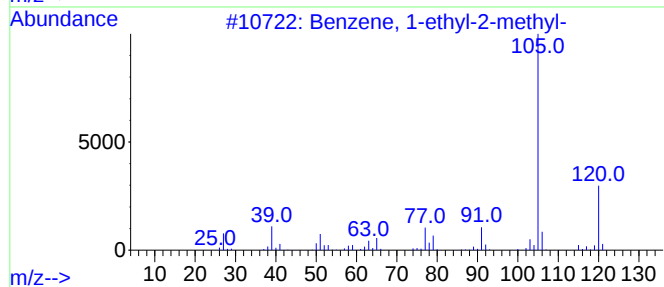
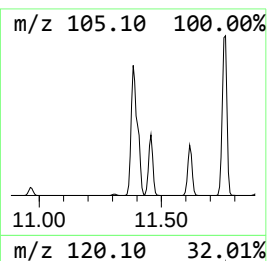
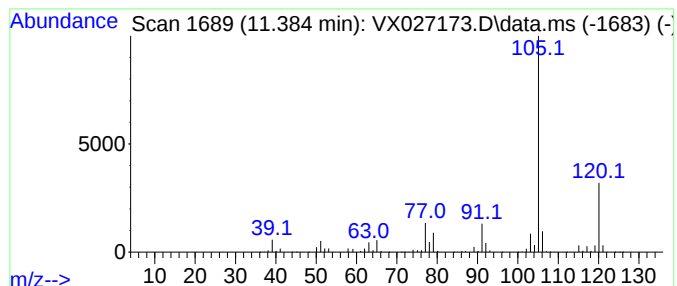
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	736.71 ug/l	8512490	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

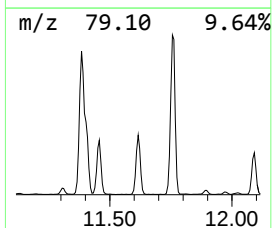
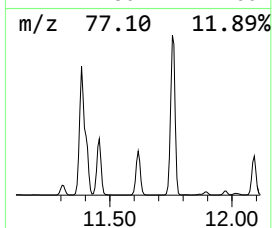
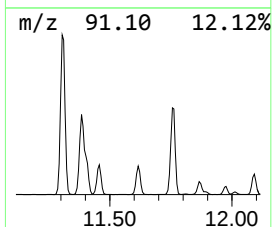
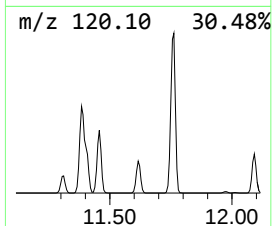
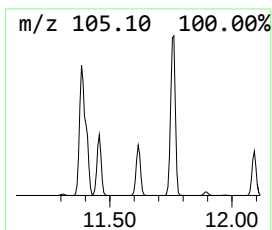
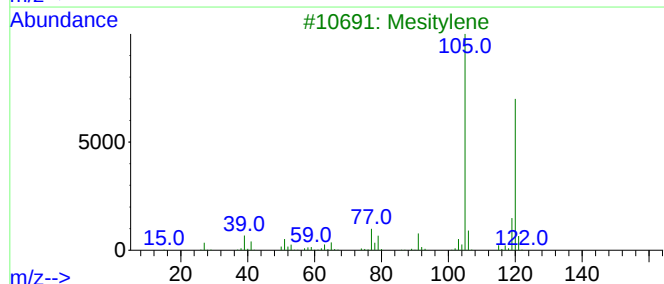
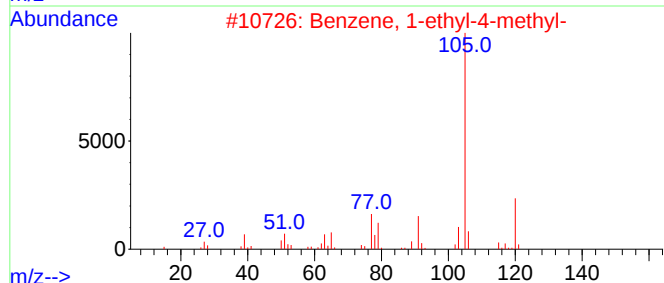
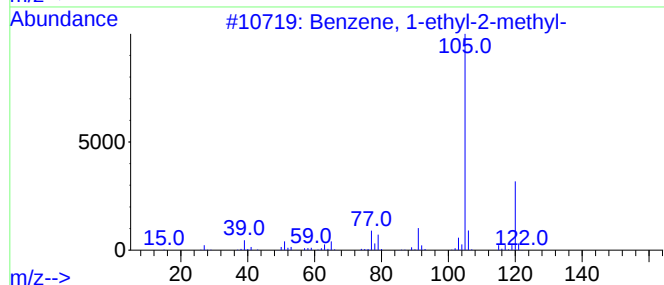
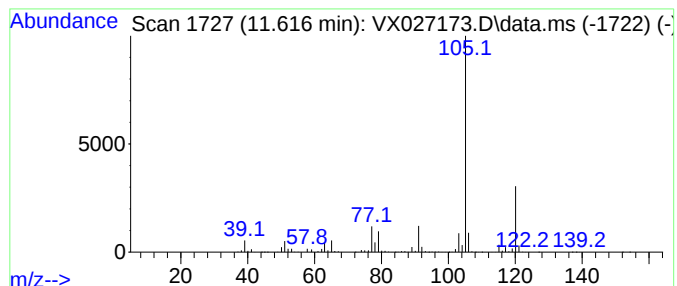
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	197.08 ug/l	2277200	1,4-Dichlorobenzene-d4	12.024

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-4-methyl-	120	C9H12	000622-96-8	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

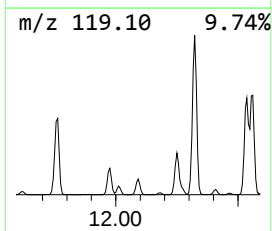
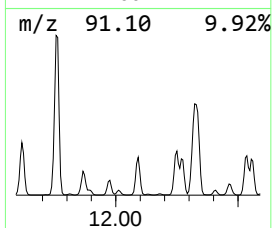
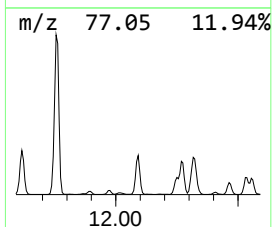
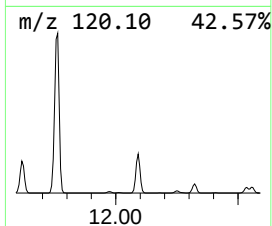
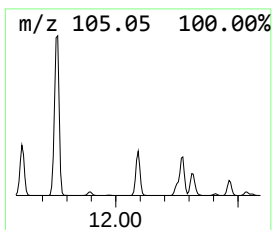
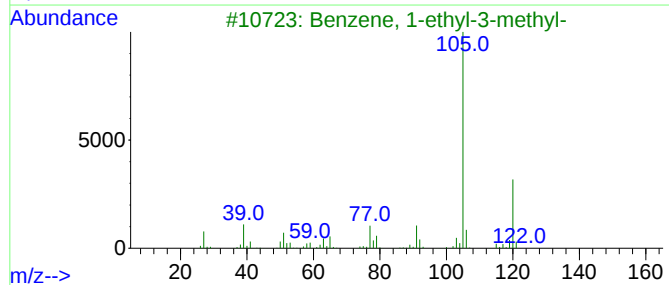
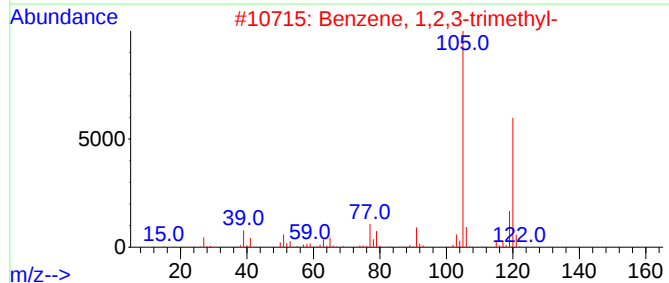
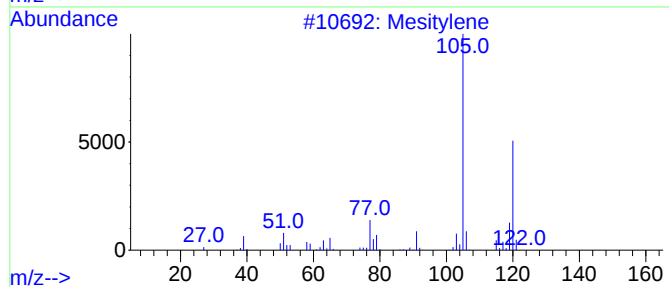
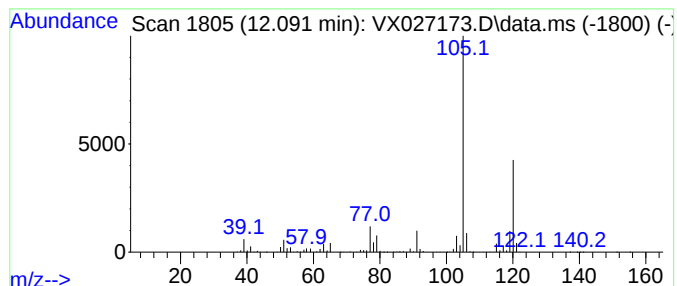
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	192.79 ug/l	2227590	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

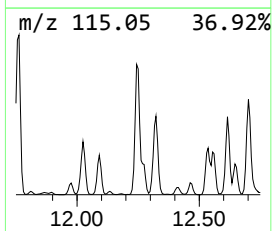
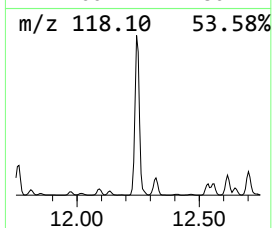
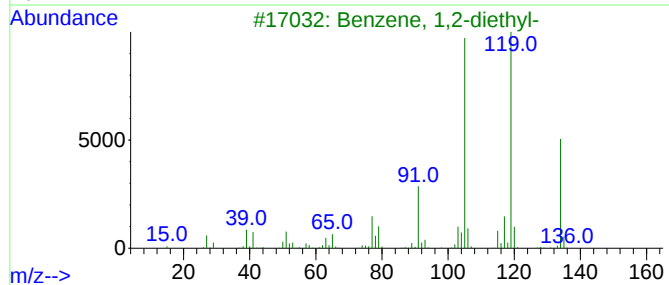
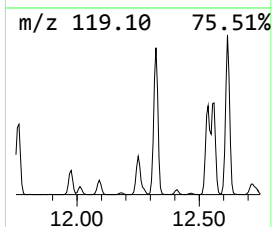
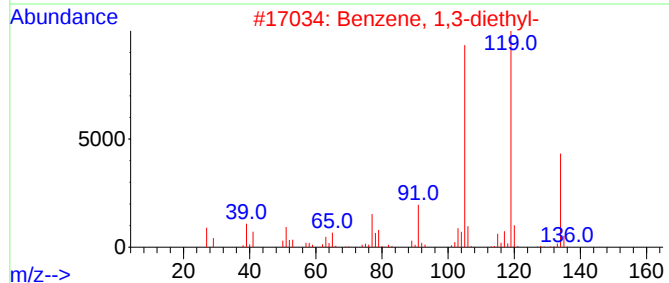
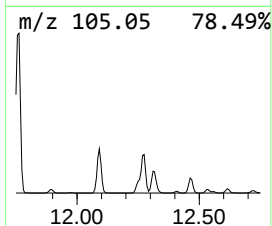
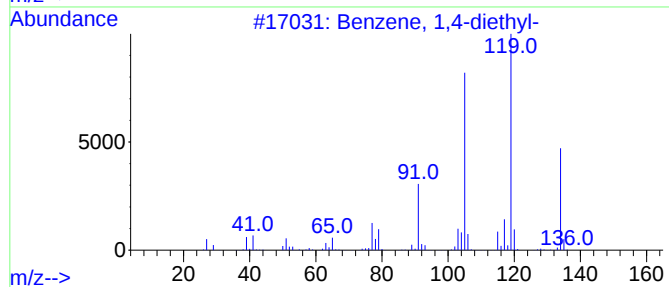
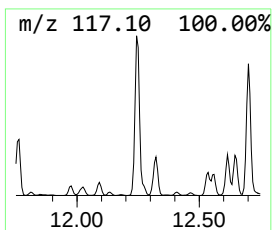
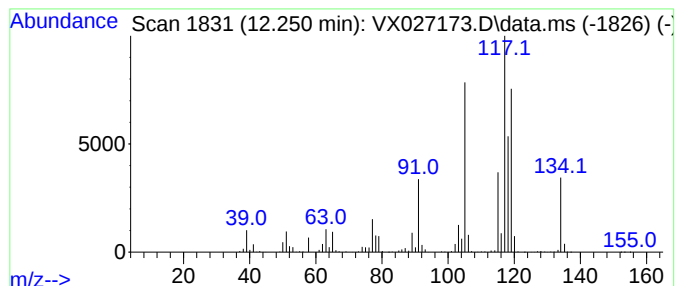
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,4-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	166.25 ug/l	1921010	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
2			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	50
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

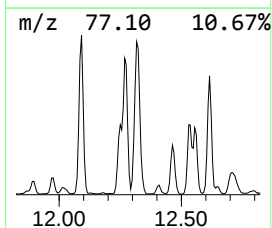
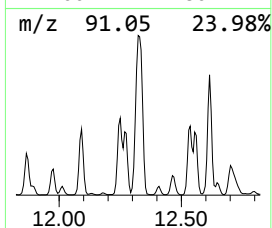
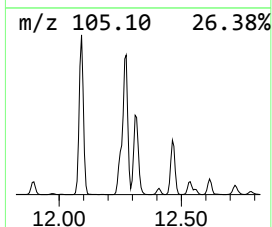
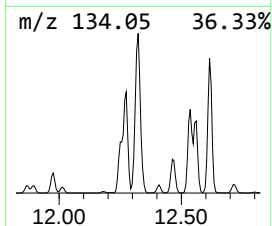
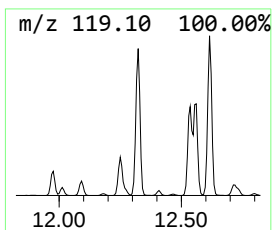
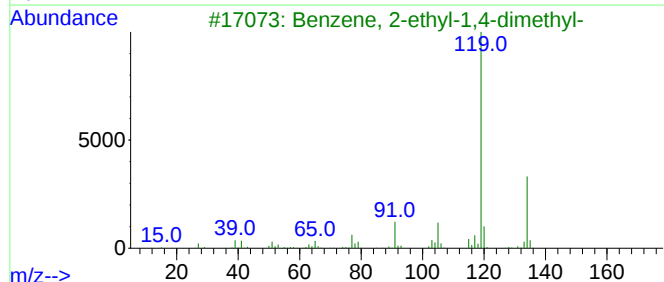
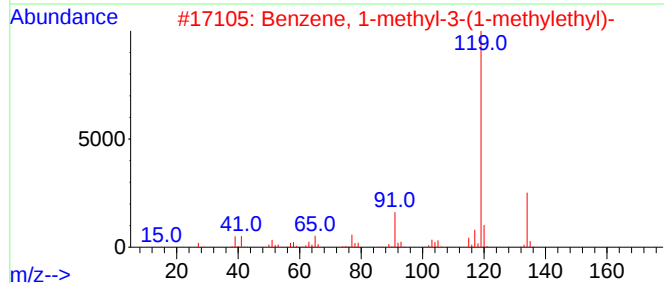
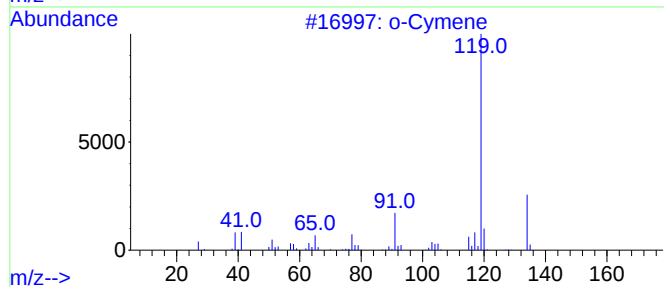
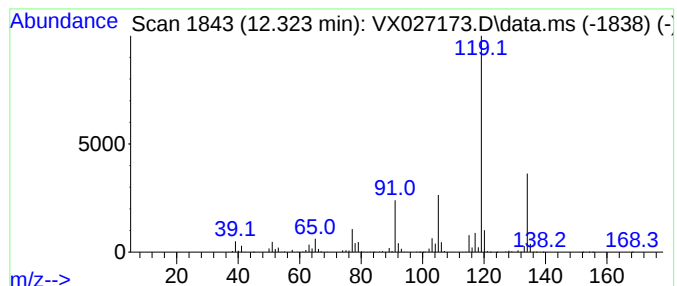
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.323	268.59 ug/l	3103550	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			p-Cymene	134	C10H14	000099-87-6	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

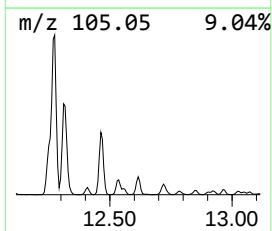
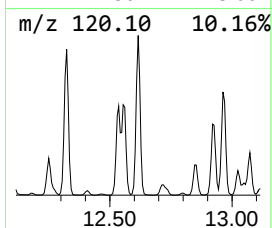
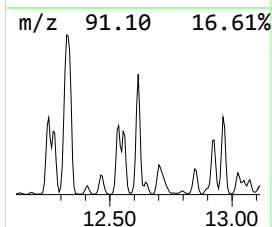
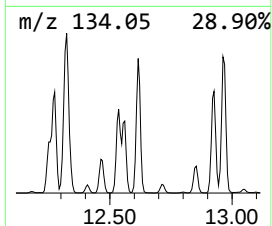
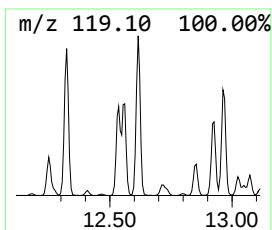
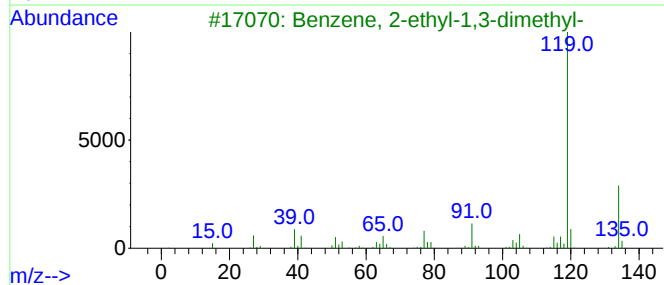
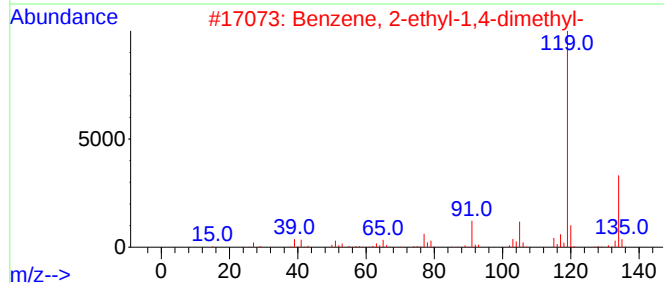
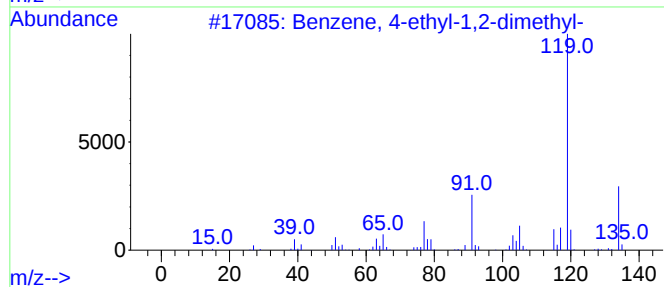
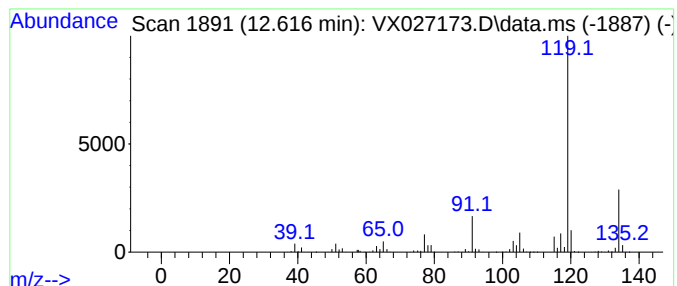
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	179.63 ug/l	2075540	1,4-Dichlorobenzene-d4	12.024

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

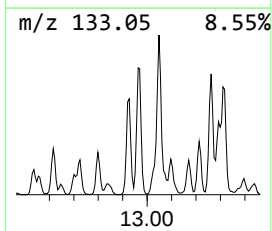
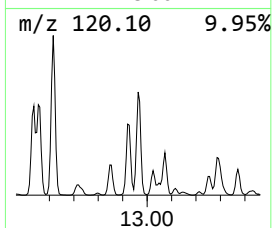
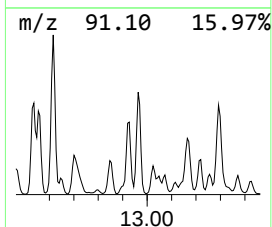
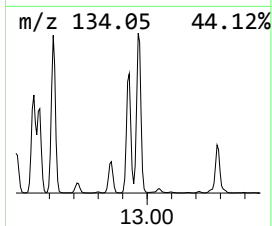
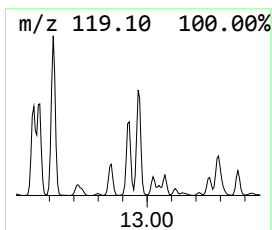
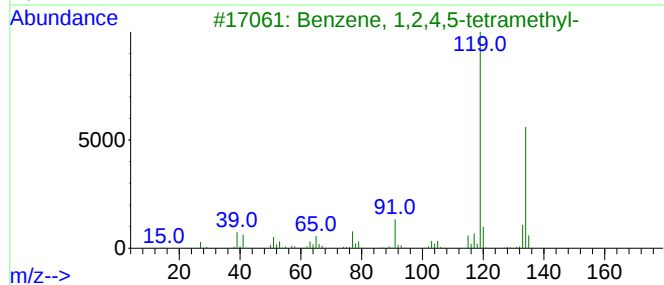
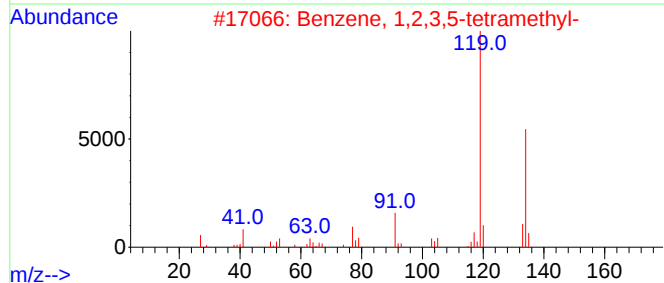
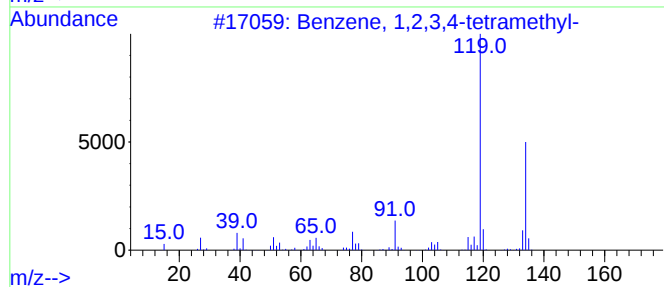
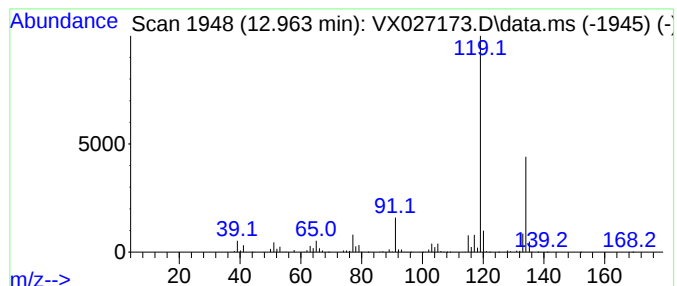
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	137.29 ug/l	1586360	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

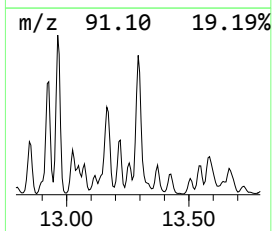
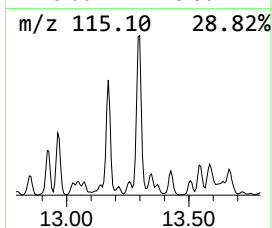
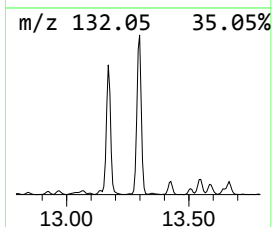
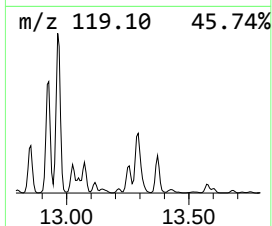
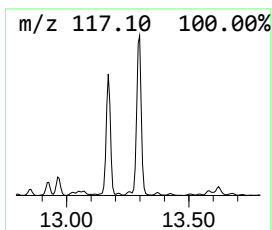
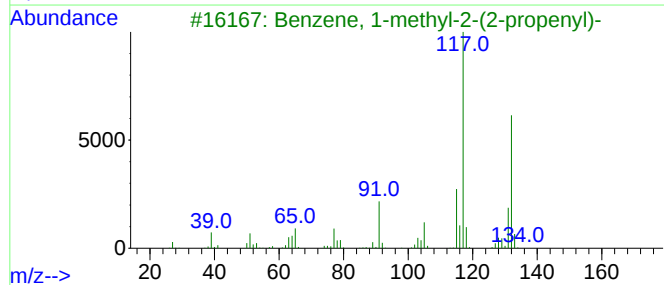
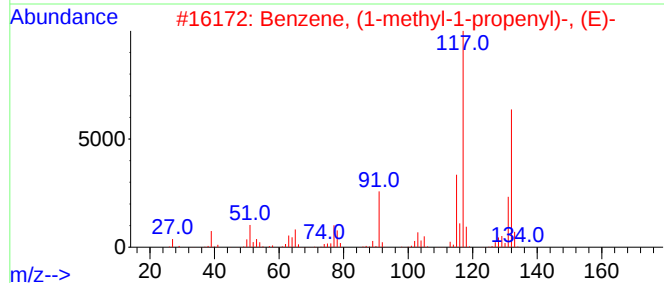
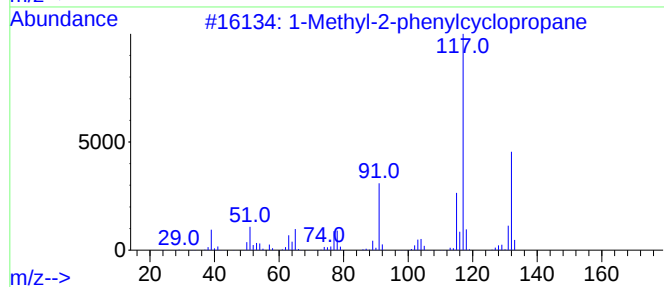
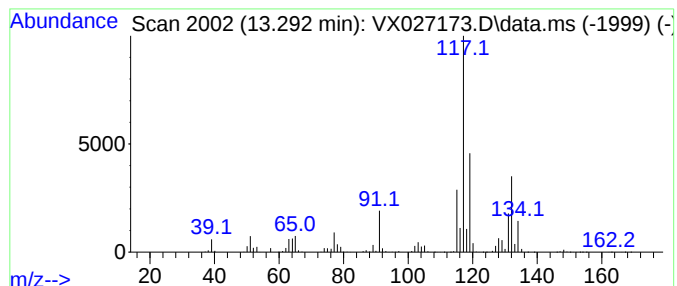
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 1-Methyl-2-phenylcyclopropane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	161.05 ug/l	1860900	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	76
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	76
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027173.D
Acq On : 24 Feb 2022 16:26
Operator : JC/MD
Sample : N1576-01
Misc : 4.93g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,2,4-...	6.245	181.7	ug/l	1273800	2	6.763	350485	50.0
Pentane, 2,3,3-...	8.104	135.9	ug/l	952604	2	6.763	350485	50.0
Benzene, 1-ethy...	11.384	736.7	ug/l	8512490	4	12.024	577740	50.0
Benzene, 1-ethy...	11.616	197.1	ug/l	2277200	4	12.024	577740	50.0
Benzene, 1,2,3-...	12.091	192.8	ug/l	2227590	4	12.024	577740	50.0
Benzene, 1,4-di...	12.250	166.3	ug/l	1921010	4	12.024	577740	50.0
o-Cymene	12.323	268.6	ug/l	3103550	4	12.024	577740	50.0
Benzene, 4-ethy...	12.616	179.6	ug/l	2075540	4	12.024	577740	50.0
Benzene, 1,2,3,...	12.963	137.3	ug/l	1586360	4	12.024	577740	50.0
1-Methyl-2-phen...	13.292	161.1	ug/l	1860900	4	12.024	577740	50.0