

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
 Data File : VX027175.D
 Acq On : 24 Feb 2022 17:13
 Operator : JC/MD
 Sample : N1539-02 10X
 Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PE-2

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: VX027175.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.825	280	285	296	rVB	73141	150161	4.19%	0.415%
2	3.099	322	330	343	rVB	55008	124588	3.48%	0.344%
3	3.446	378	387	398	rBV	40837	92443	2.58%	0.255%
4	4.215	503	513	520	rVV	26492	75991	2.12%	0.210%
5	4.306	520	528	539	rVB	57272	161126	4.50%	0.445%
6	5.196	665	674	689	rVB2	19772	68095	1.90%	0.188%
7	5.391	691	706	714	rBV	54547	171829	4.79%	0.475%
8	5.507	714	725	729	rBV3	91255	279025	7.79%	0.771%
9	5.623	739	744	765	rVB2	71258	236518	6.60%	0.654%
10	5.830	765	778	790	rBV	110858	327703	9.14%	0.906%
11	5.958	790	799	812	rVB	54830	145395	4.06%	0.402%
12	6.159	817	832	838	rBV3	39957	132668	3.70%	0.367%
13	6.251	838	847	858	rVV	167681	524543	14.64%	1.450%
14	6.361	858	865	875	rVB3	31949	91801	2.56%	0.254%
15	6.611	896	906	914	rBV2	75648	184920	5.16%	0.511%
16	6.769	924	932	938	rBV	137773	344464	9.61%	0.952%
17	6.848	939	945	952	rVV4	53909	162443	4.53%	0.449%
18	7.177	989	999	1007	rBV3	31616	79456	2.22%	0.220%
19	7.379	1018	1032	1044	rBV3	123611	344711	9.62%	0.953%
20	7.501	1044	1052	1057	rBV2	48132	99407	2.77%	0.275%
21	7.562	1057	1062	1073	rVB	55217	118557	3.31%	0.328%
22	7.659	1073	1078	1089	rVB	42214	79187	2.21%	0.219%
23	7.775	1089	1097	1105	rVB	32999	66613	1.86%	0.184%
24	7.988	1121	1132	1143	rVB	124007	288191	8.04%	0.796%
25	8.104	1143	1151	1161	rBV2	157822	409147	11.42%	1.131%
26	8.190	1161	1165	1174	rVB3	64995	123977	3.46%	0.343%
27	8.275	1174	1179	1183	rBV	130779	219461	6.12%	0.607%
28	8.311	1183	1185	1191	rVB2	66223	97293	2.71%	0.269%
29	8.439	1201	1206	1214	rBV	145465	265368	7.40%	0.733%
30	8.507	1214	1217	1227	rVB4	39153	85937	2.40%	0.238%
31	8.610	1227	1234	1236	rBV3	94534	180830	5.05%	0.500%
32	8.653	1236	1241	1246	rVB	295187	500556	13.97%	1.383%
33	8.720	1246	1252	1258	rBV	1466515	2209570	61.66%	6.107%
34	8.915	1277	1284	1291	rBV3	114705	234350	6.54%	0.648%
35	8.982	1291	1295	1305	rVB5	31270	79916	2.23%	0.221%
36	9.104	1311	1315	1321	rVB2	44019	65737	1.83%	0.182%

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37	9.159	1321	1324	1328	rVV2	36752	54637	1.52%	0.151%
38	9.202	1328	1331	1341	rVB4	24865	52306	1.46%	0.145%
39	9.384	1355	1361	1366	rBV2	46061	74646	2.08%	0.206%
40	9.500	1375	1380	1387	rBV2	81454	163331	4.56%	0.451%
41	9.653	1403	1405	1416	rVB8	21365	56659	1.58%	0.157%
42	9.762	1416	1423	1428	rBV5	26915	54940	1.53%	0.152%
43	9.823	1428	1433	1437	rBV3	30675	61932	1.73%	0.171%
44	9.903	1442	1446	1452	rVB	188775	270590	7.55%	0.748%
45	10.012	1458	1464	1467	rBV	143903	195170	5.45%	0.539%
46	10.055	1467	1471	1476	rVB	311686	395674	11.04%	1.094%
47	10.195	1489	1494	1500	rVB	945111	1200852	33.51%	3.319%
48	10.305	1500	1512	1518	rVV	2588469	3583753	100.00%	9.904%
49	10.360	1518	1521	1532	rVB2	140651	235730	6.58%	0.651%
50	10.585	1553	1558	1562	rBV	58672	82729	2.31%	0.229%
51	10.640	1562	1567	1577	rVV	1509827	2078928	58.01%	5.746%
52	10.781	1582	1590	1594	rBV	62569	101308	2.83%	0.280%
53	10.860	1594	1603	1607	rBV3	49007	81754	2.28%	0.226%
54	10.963	1614	1620	1624	rBV	115173	152653	4.26%	0.422%
55	11.030	1627	1631	1634	rVV2	39192	55018	1.54%	0.152%
56	11.085	1634	1640	1643	rVV	327969	497747	13.89%	1.376%
57	11.110	1643	1644	1651	rVV2	125320	152654	4.26%	0.422%
58	11.195	1651	1658	1665	rVV	98433	146448	4.09%	0.405%
59	11.305	1672	1676	1683	rVB	366245	437977	12.22%	1.210%
60	11.384	1683	1689	1696	rBV	1647074	2906572	81.10%	8.033%
61	11.451	1696	1700	1704	rBV	811417	1014198	28.30%	2.803%
62	11.616	1722	1727	1734	rVB	619063	803638	22.42%	2.221%
63	11.756	1745	1750	1760	rVB	2536161	3112514	86.85%	8.602%
64	11.890	1770	1772	1776	rVB	53545	59943	1.67%	0.166%
65	11.969	1782	1785	1789	rVB	84591	97183	2.71%	0.269%
66	12.024	1789	1794	1799	rBV	340772	475903	13.28%	1.315%
67	12.091	1799	1805	1810	rBV	553980	772766	21.56%	2.136%
68	12.177	1816	1819	1823	rVB2	62258	69844	1.95%	0.193%
69	12.244	1825	1830	1832	rBV2	472362	608204	16.97%	1.681%
70	12.268	1832	1834	1838	rVV	426000	475150	13.26%	1.313%
71	12.317	1838	1842	1849	rVB3	587609	916515	25.57%	2.533%
72	12.420	1849	1859	1863	rBV3	123454	227474	6.35%	0.629%
73	12.463	1863	1866	1872	rVB	153891	196857	5.49%	0.544%
74	12.530	1872	1877	1879	rBV	294702	372262	10.39%	1.029%
75	12.555	1879	1881	1887	rVV	294335	357731	9.98%	0.989%
76	12.616	1887	1891	1894	rVV	482284	572480	15.97%	1.582%

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77	12.646	1894	1896	1900	rVV	69006	72124	2.01%	0.199%
78	12.701	1900	1905	1916	rVB3	159815	306351	8.55%	0.847%
79	12.847	1925	1929	1933	rVB	95352	111789	3.12%	0.309%
80	12.920	1933	1941	1945	rBV	252191	373458	10.42%	1.032%
81	12.963	1945	1948	1954	rVB	350231	449970	12.56%	1.244%
82	13.024	1954	1958	1959	rBV	69366	75063	2.09%	0.207%
83	13.042	1959	1961	1964	rVV3	84294	117053	3.27%	0.324%
84	13.073	1964	1966	1969	rVB	55972	53093	1.48%	0.147%
85	13.170	1974	1982	1986	rVV3	215672	325891	9.09%	0.901%
86	13.213	1986	1989	1992	rVV2	51725	56870	1.59%	0.157%
87	13.268	1992	1998	1999	rVV3	101635	168794	4.71%	0.466%
88	13.292	1999	2002	2008	rVV2	377297	526112	14.68%	1.454%
89	13.372	2012	2015	2020	rVB	58249	71470	1.99%	0.198%
90	13.426	2020	2024	2032	rVB4	47835	88373	2.47%	0.244%
91	13.506	2032	2037	2040	rBV2	49595	71055	1.98%	0.196%
92	13.542	2040	2043	2046	rVV	65529	84382	2.35%	0.233%
93	13.585	2046	2050	2056	rVV2	91984	173604	4.84%	0.480%
94	13.664	2060	2063	2069	rVB2	55788	95632	2.67%	0.264%
95	13.774	2077	2081	2088	rVB	304114	384557	10.73%	1.063%
96	13.926	2100	2106	2111	rBV3	38406	71722	2.00%	0.198%
97	14.079	2127	2131	2137	rVB	47155	59307	1.65%	0.164%
98	14.213	2149	2153	2160	rBV	39840	63212	1.76%	0.175%
99	14.640	2214	2223	2233	rVB	166703	233155	6.51%	0.644%
100	14.780	2240	2246	2255	rVB	75498	101618	2.84%	0.281%

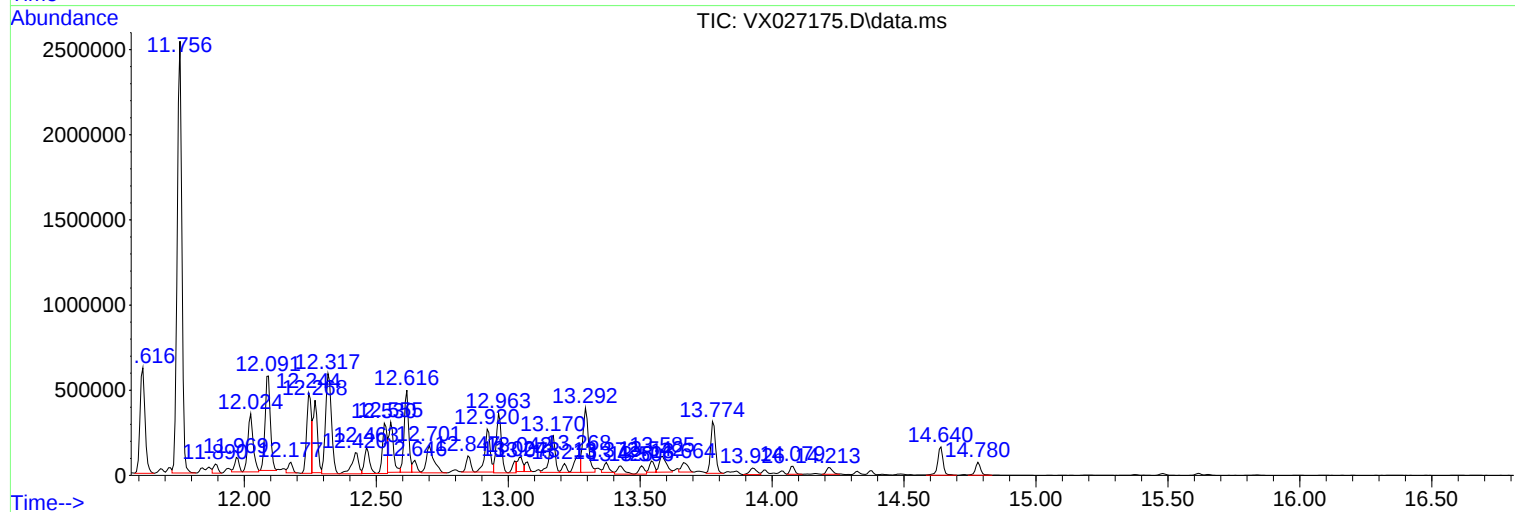
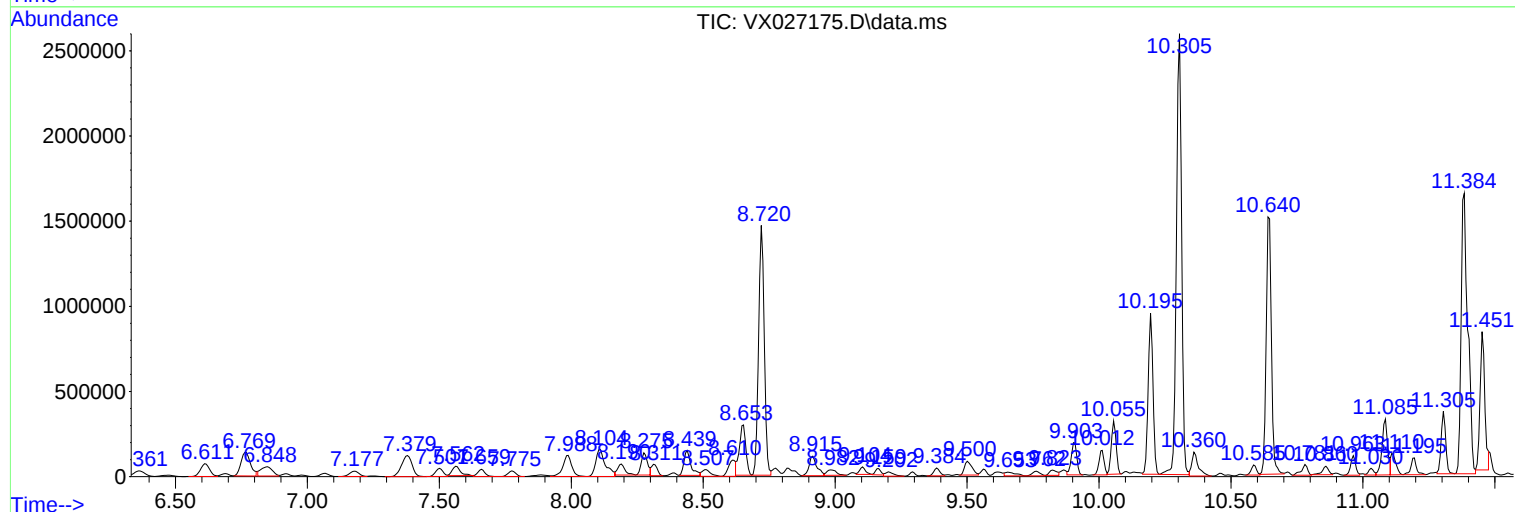
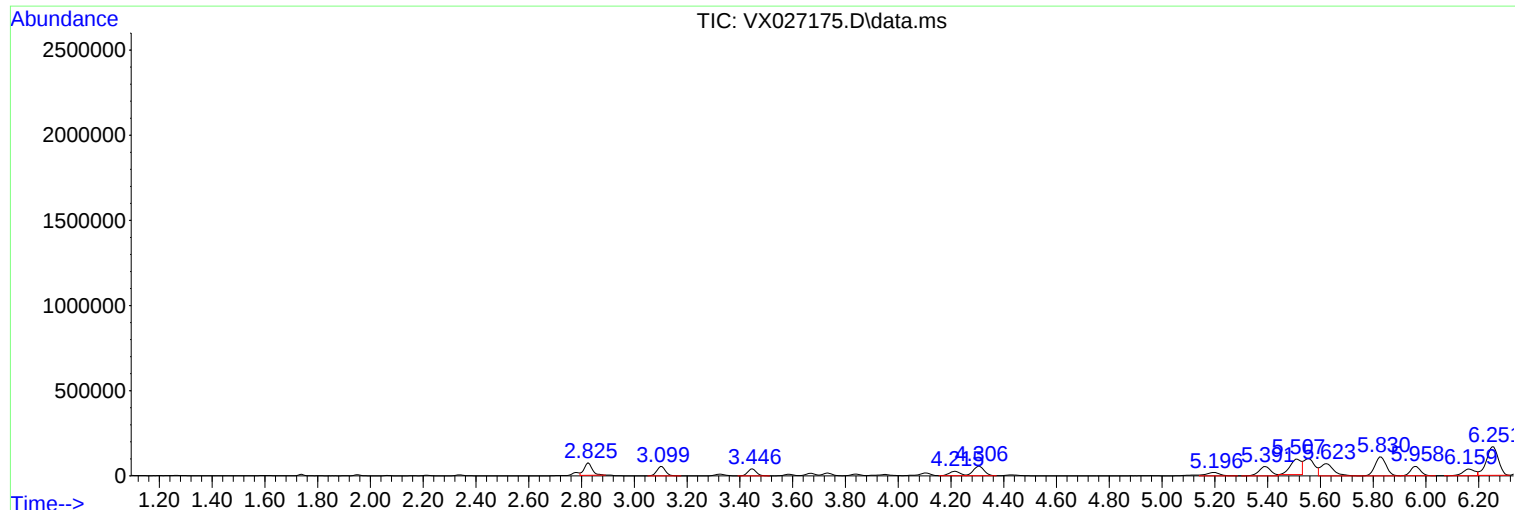
Sum of corrected areas: 36183302

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Instrument :
MSVOA_X
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PE-2

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



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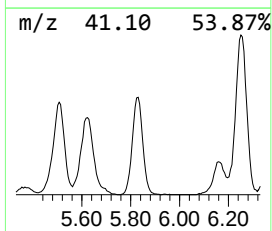
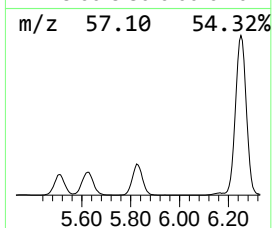
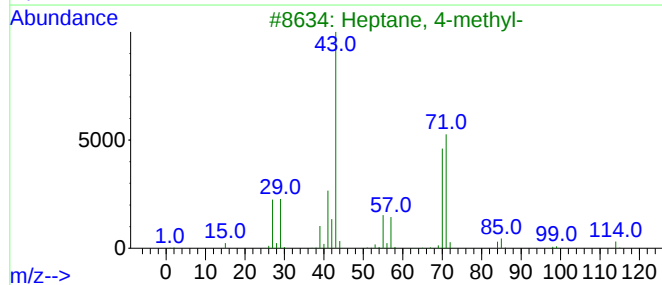
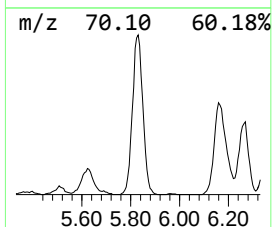
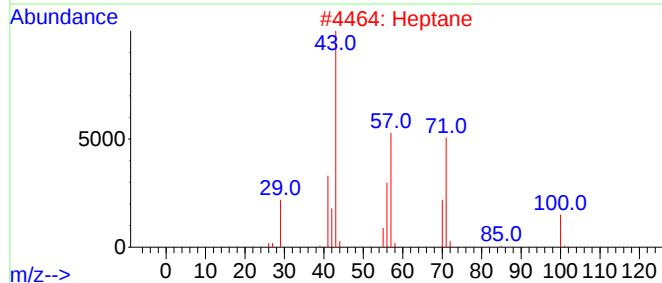
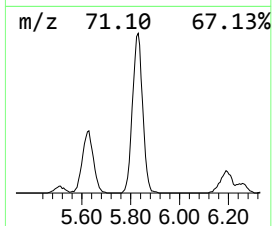
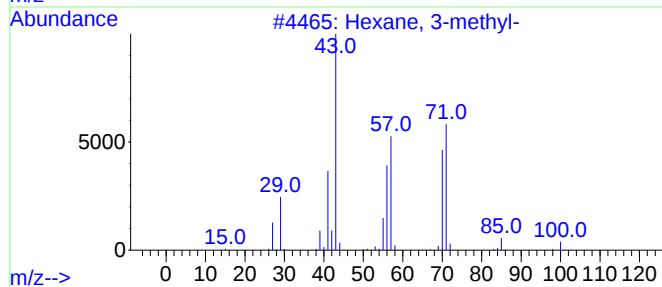
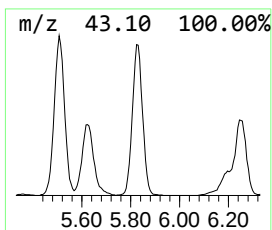
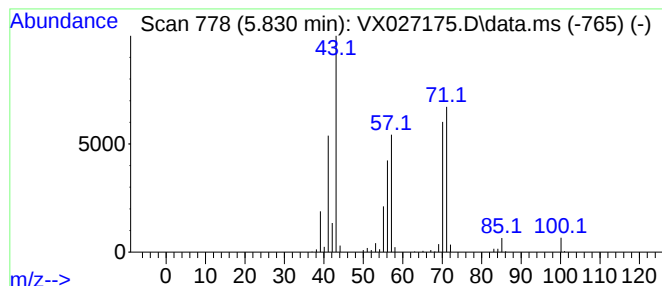
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022322W.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.830	58.72 ug/l	327703	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53



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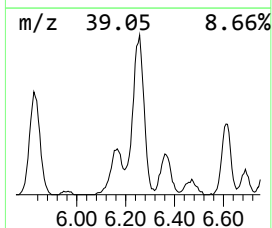
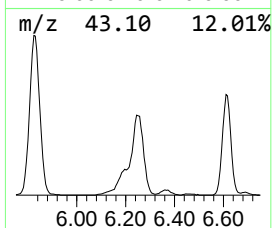
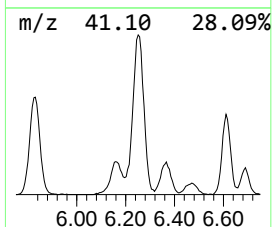
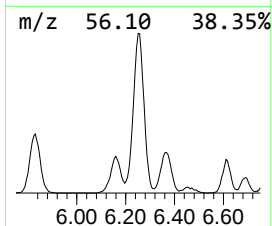
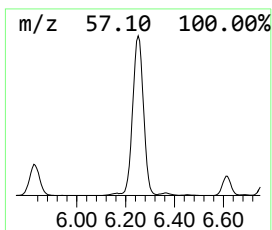
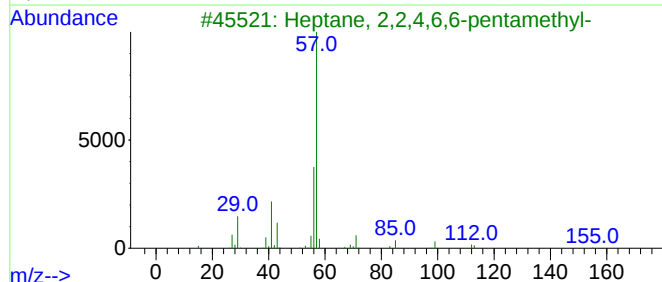
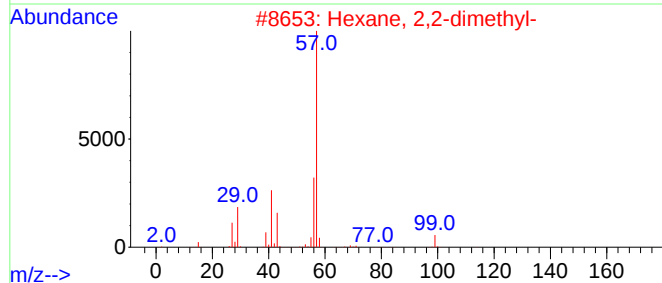
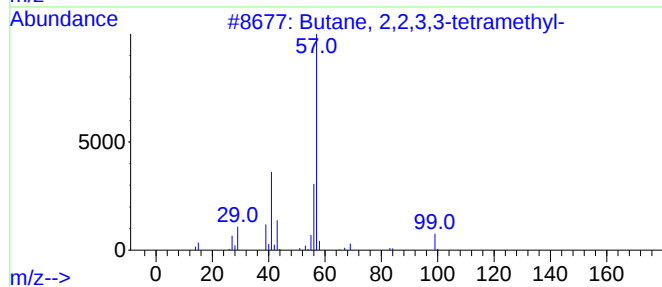
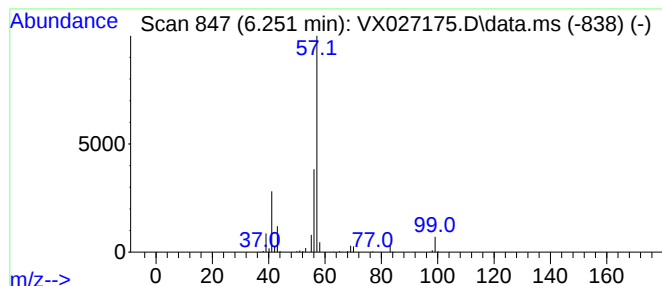
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Quant Title : SW846 8260

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

Peak Number 2 Butane, 2,2,3,3-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.251	76.14 ug/l	524543	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	72
4			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
5			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64



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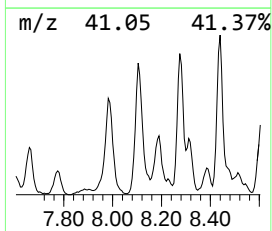
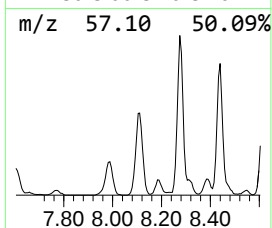
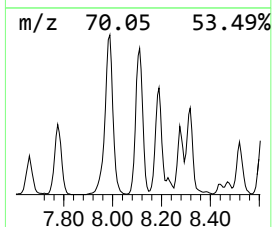
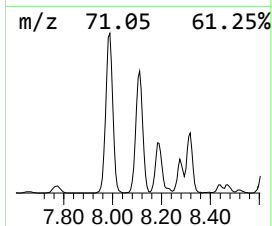
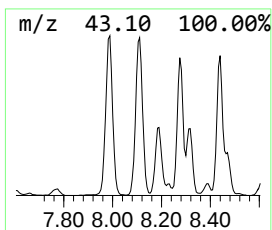
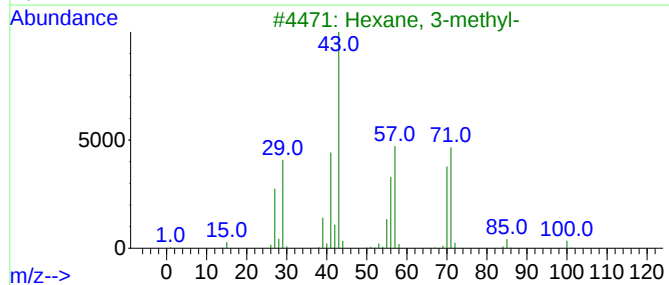
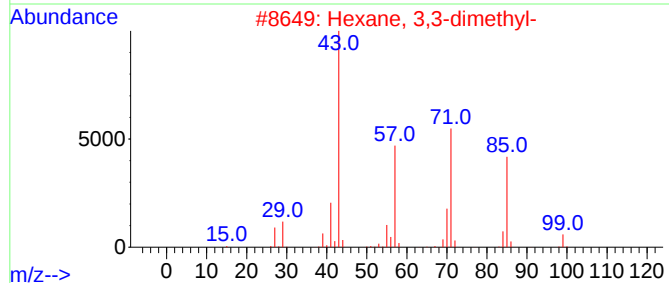
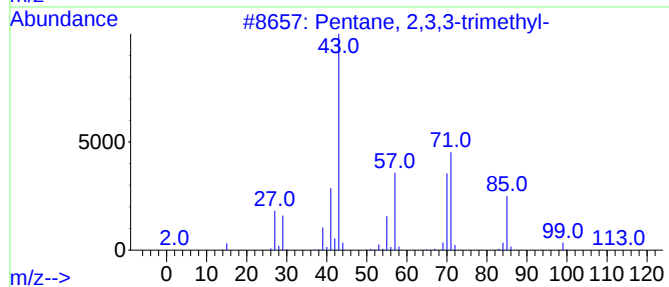
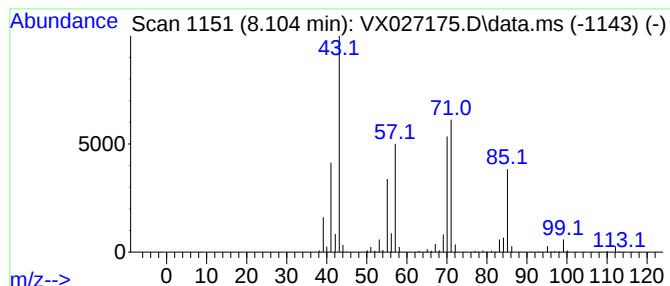
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	59.39 ug/l	409147	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			Hexane, 3-methyl-	100	C7H16	000589-34-4	58
4			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	53
5			Heptane, 4-methyl-	114	C8H18	000589-53-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

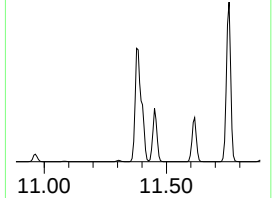
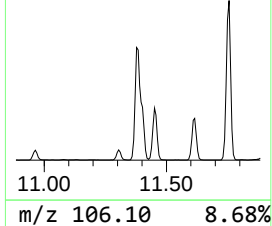
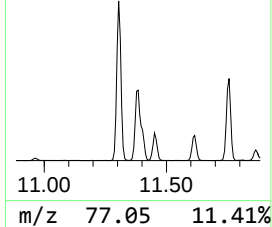
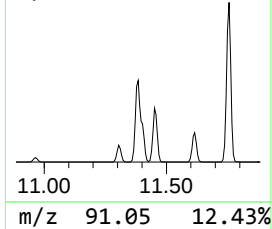
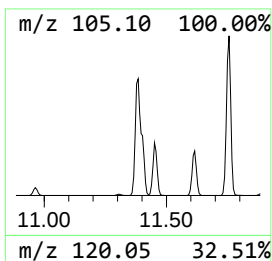
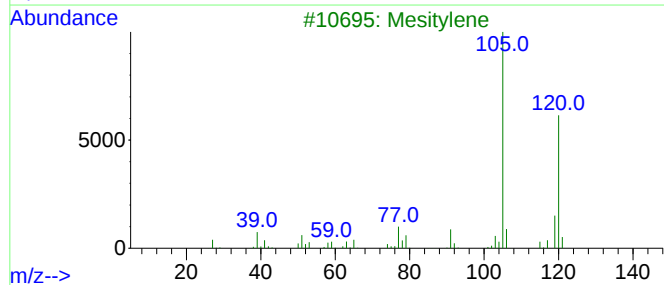
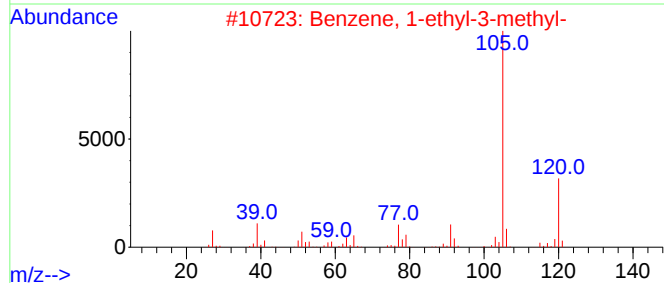
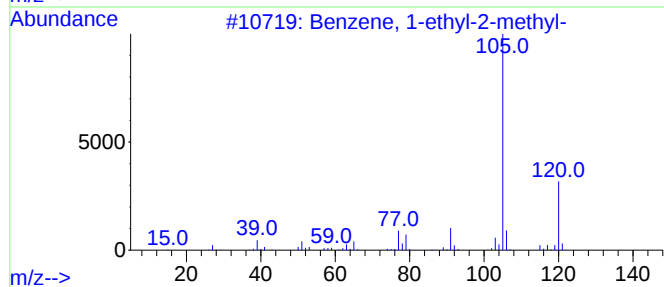
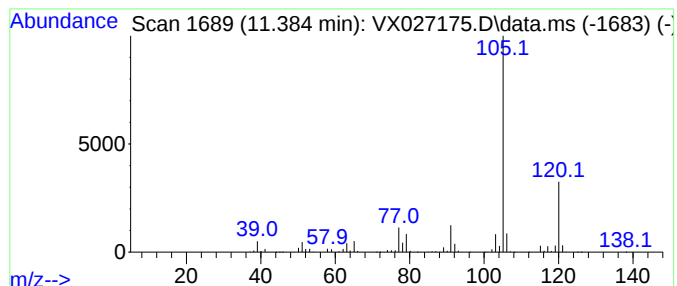
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	305.37 ug/l	2906570	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			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

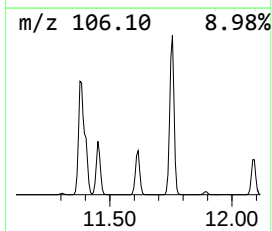
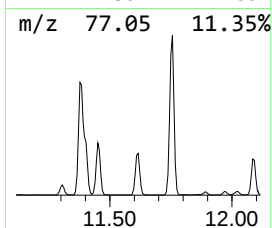
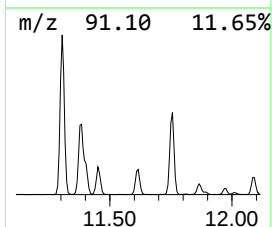
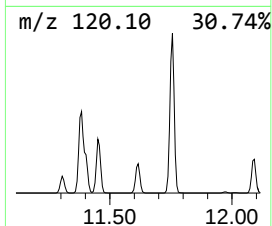
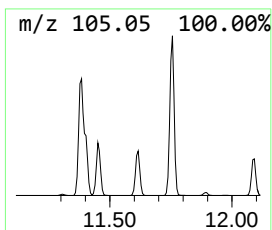
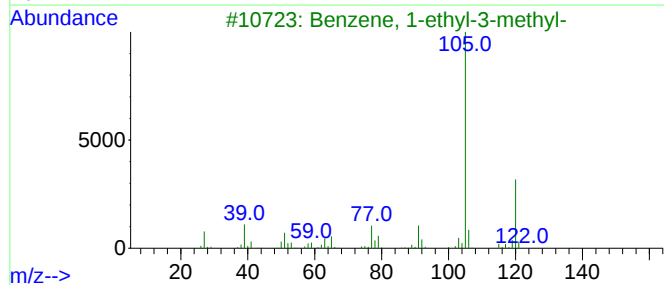
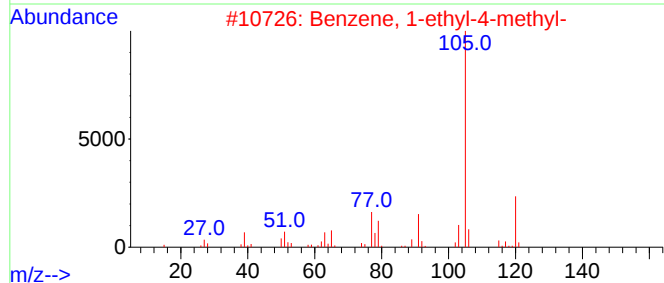
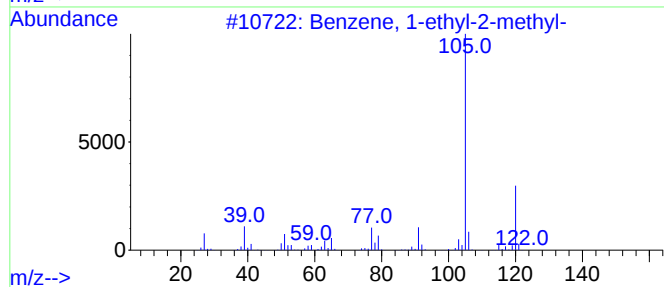
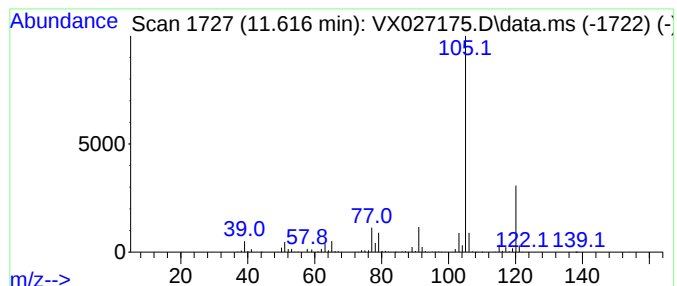
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	84.43 ug/l	803638	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-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

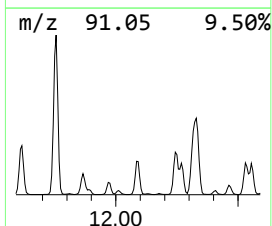
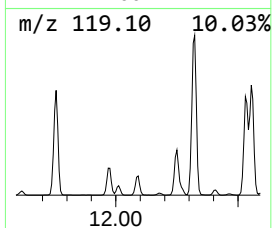
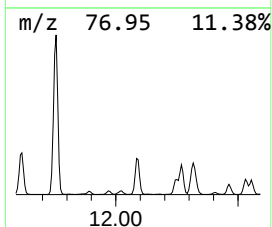
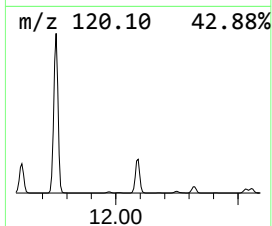
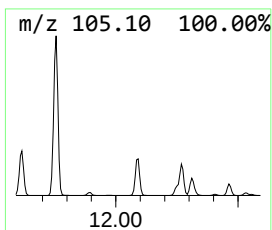
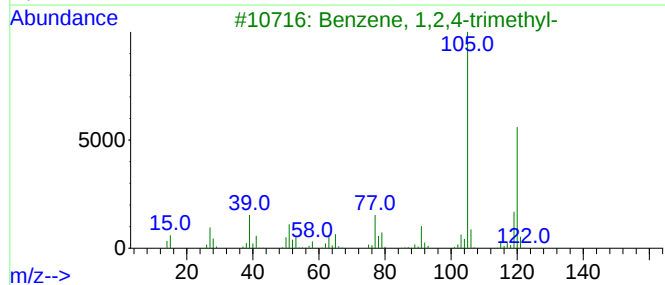
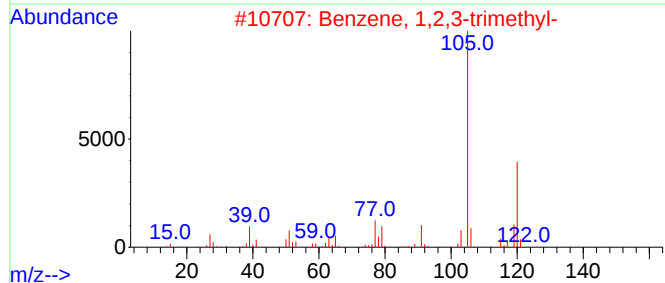
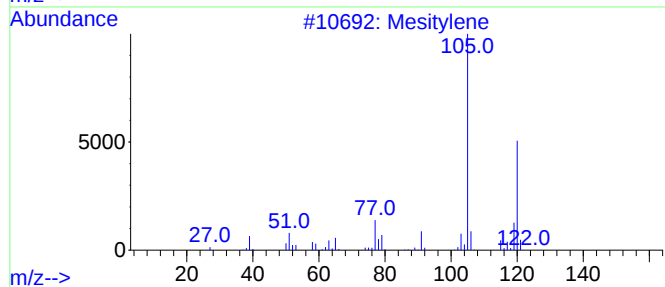
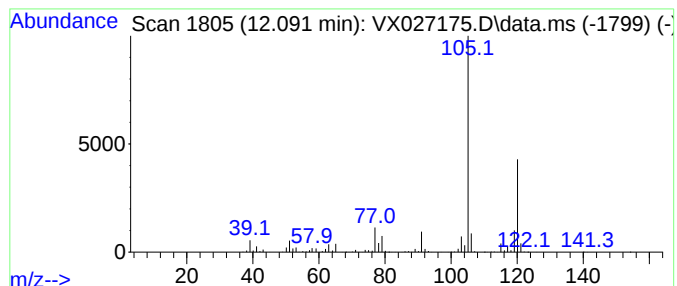
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	81.19 ug/l	772766	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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

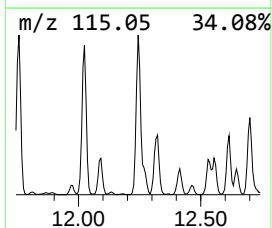
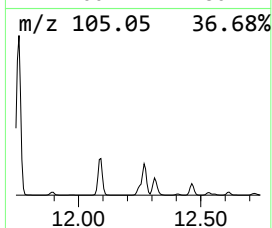
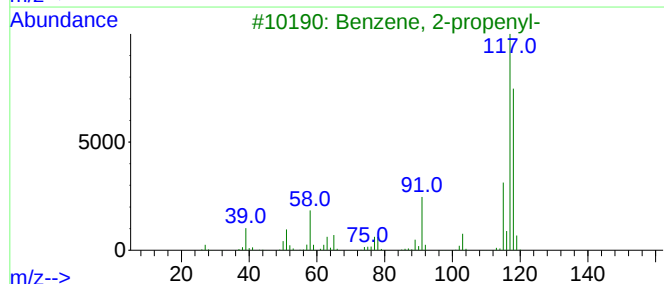
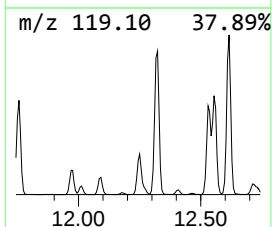
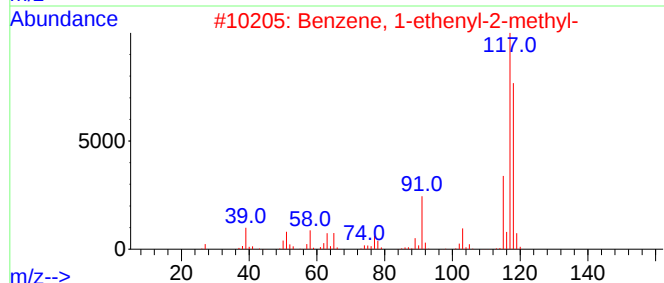
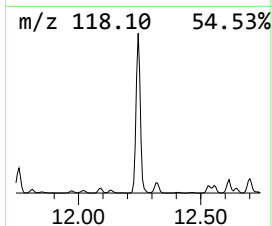
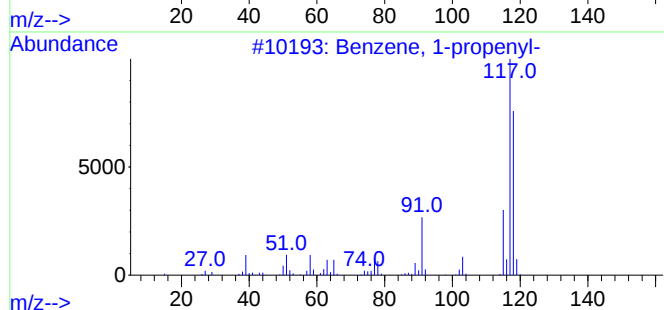
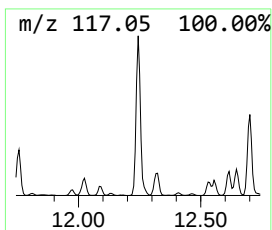
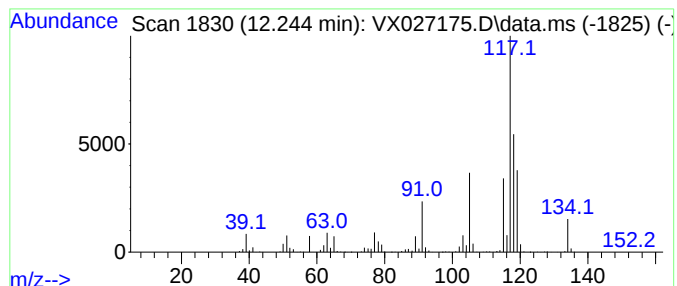
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	63.90 ug/l	608204	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	92
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			Indane	118	C9H10	000496-11-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

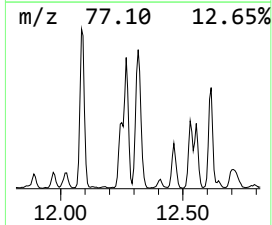
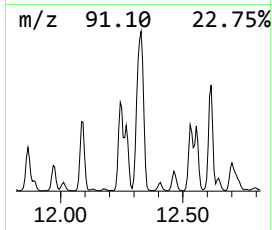
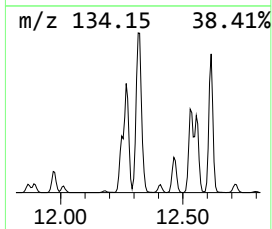
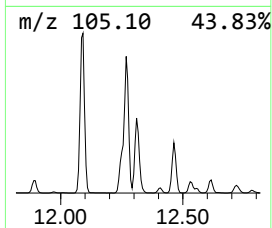
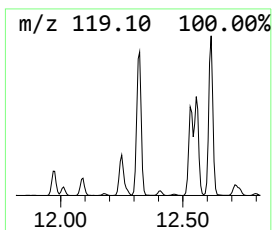
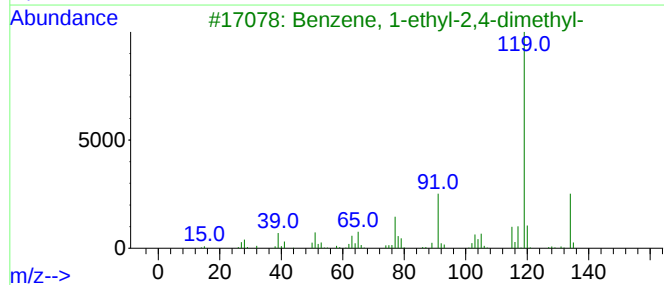
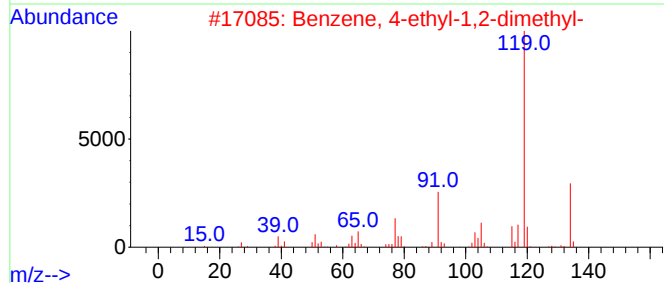
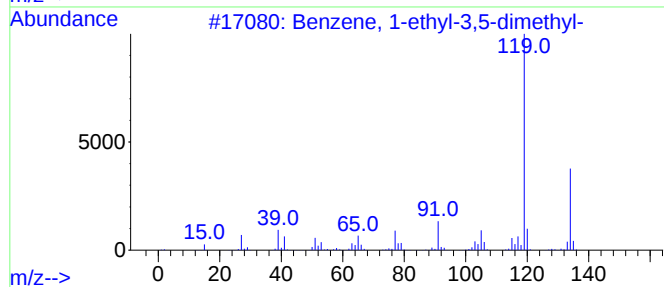
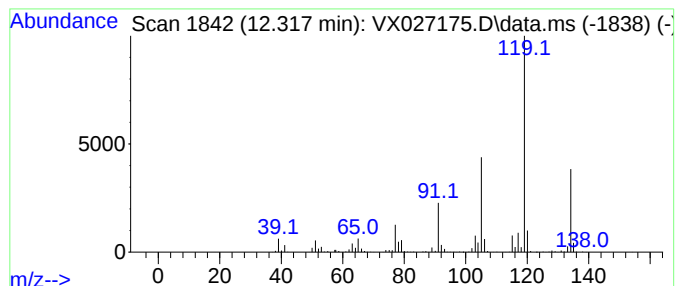
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	96.29 ug/l	916515	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

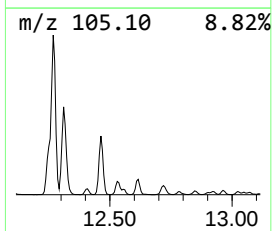
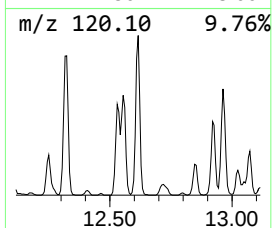
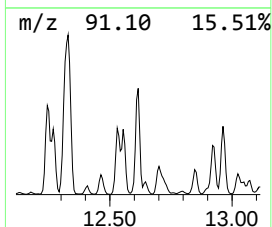
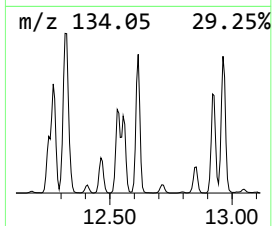
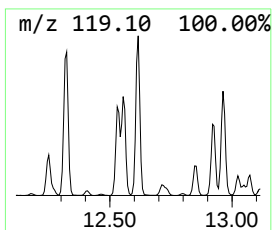
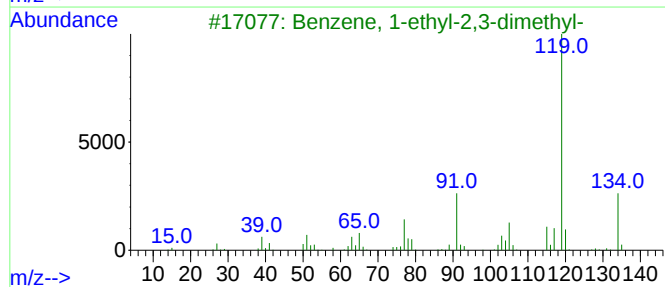
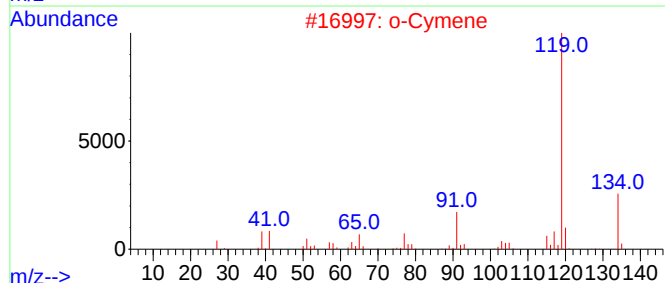
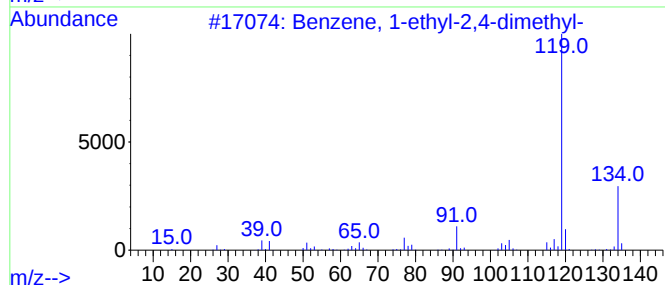
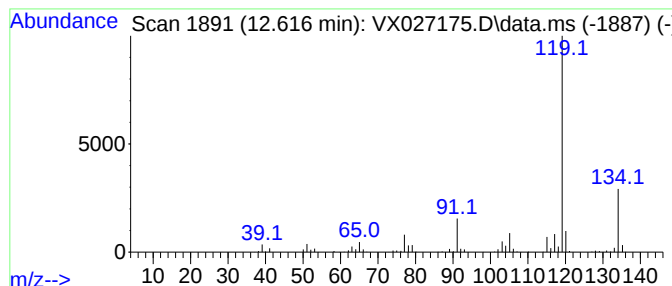
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	60.15 ug/l	572480	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

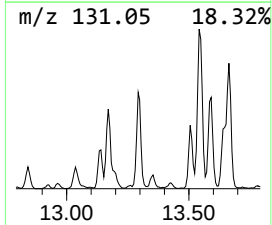
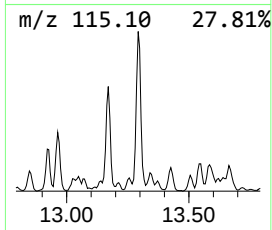
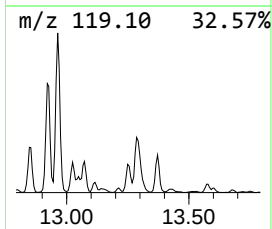
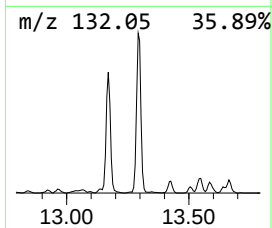
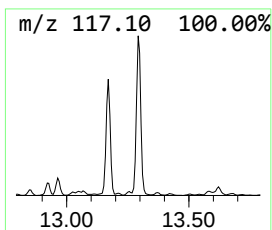
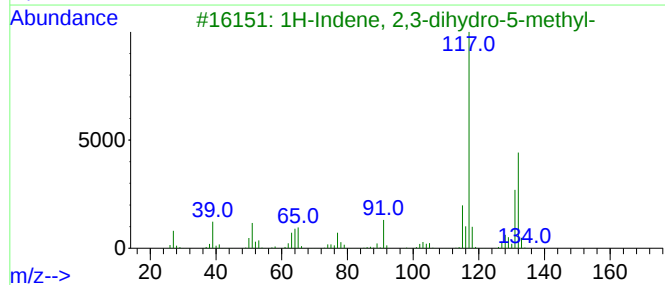
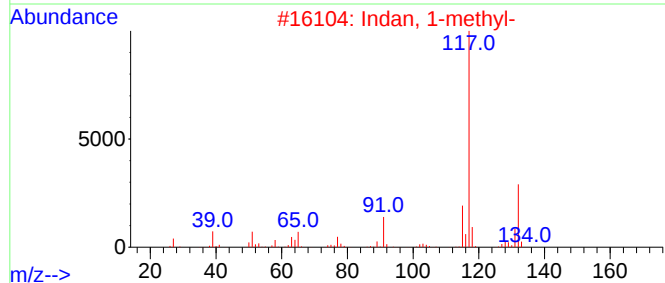
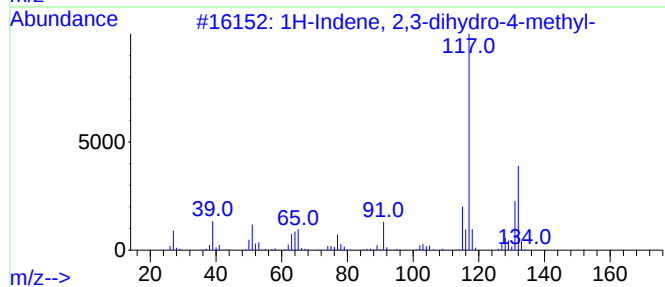
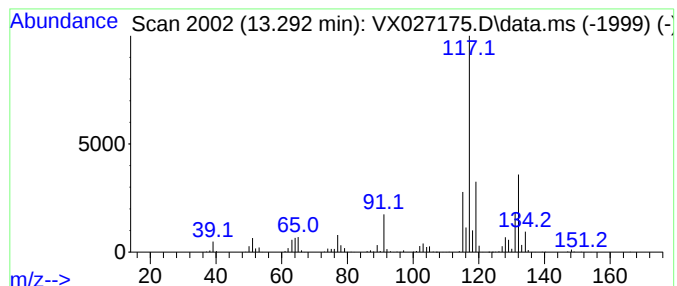
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	55.28 ug/l	526112	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2			Indan, 1-methyl-	132	C10H12	000767-58-8	93
3			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	81
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022422\
Data File : VX027175.D
Acq On : 24 Feb 2022 17:13
Operator : JC/MD
Sample : N1539-02 10X
Misc : 5.31g/5.0mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PE-2

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
Hexane, 3-methyl-	5.830	58.7	ug/l	327703	1	5.556	279025	50.0
Butane, 2,2,3,3...	6.251	76.1	ug/l	524543	2	6.763	344464	50.0
Pentane, 2,3,3-...	8.104	59.4	ug/l	409147	2	6.763	344464	50.0
Benzene, 1-ethy...	11.384	305.4	ug/l	2906570	4	12.024	475903	50.0
Benzene, 1-ethy...	11.616	84.4	ug/l	803638	4	12.024	475903	50.0
Benzene, 1,2,3-...	12.091	81.2	ug/l	772766	4	12.024	475903	50.0
Benzene, 1-prop...	12.244	63.9	ug/l	608204	4	12.024	475903	50.0
Benzene, 1-ethy...	12.317	96.3	ug/l	916515	4	12.024	475903	50.0
Benzene, 1-ethy...	12.616	60.1	ug/l	572480	4	12.024	475903	50.0
1H-Indene, 2,3-...	13.292	55.3	ug/l	526112	4	12.024	475903	50.0