

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
 Data File : VX045822.D
 Acq On : 17 Apr 2025 12:49
 Operator : JC/MD
 Sample : Q1824-04 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 285

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\82X040225W.M
 Title : SW846 8260

Signal : TIC: VX045822.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.379	694	704	713	rBV	52588	154979	1.60%	0.290%
2	5.550	713	732	740	rVV3	72365	301591	3.12%	0.565%
3	5.818	762	776	789	rBV	45849	139897	1.45%	0.262%
4	5.952	789	798	805	rBV2	58916	159747	1.65%	0.299%
5	6.031	805	811	823	rVB	100781	265994	2.75%	0.498%
6	6.239	837	845	856	rVB3	57333	180298	1.86%	0.338%
7	6.604	896	905	914	rBV	71265	177445	1.83%	0.332%
8	6.757	922	930	939	rBV	131537	321762	3.32%	0.603%
9	7.373	1018	1031	1042	rBV2	143884	363196	3.75%	0.680%
10	7.976	1120	1130	1143	rVB2	67834	170301	1.76%	0.319%
11	8.104	1143	1151	1155	rBV	73169	160248	1.66%	0.300%
12	8.269	1173	1178	1182	rBV	134782	224828	2.32%	0.421%
13	8.433	1200	1205	1213	rBV	134187	237721	2.46%	0.445%
14	8.598	1225	1232	1236	rBV4	115524	260187	2.69%	0.487%
15	8.647	1236	1240	1245	rVB	274543	444983	4.60%	0.833%
16	8.714	1245	1251	1258	rBV	4976605	8076317	83.43%	15.126%
17	8.909	1277	1283	1290	rBV	199647	317060	3.28%	0.594%
18	8.976	1290	1294	1305	rVB2	53388	105190	1.09%	0.197%
19	9.098	1305	1314	1320	rBV	82851	155325	1.60%	0.291%
20	9.378	1354	1360	1370	rBV2	54592	101156	1.04%	0.189%
21	9.500	1375	1380	1386	rBV3	89260	194928	2.01%	0.365%
22	9.561	1386	1390	1394	rVB	91541	127521	1.32%	0.239%
23	9.823	1427	1433	1437	rBV3	45599	98547	1.02%	0.185%
24	9.903	1441	1446	1451	rVB	178262	259861	2.68%	0.487%
25	10.006	1457	1463	1467	rBV	139445	189778	1.96%	0.355%
26	10.049	1467	1470	1476	rVB	253347	344789	3.56%	0.646%
27	10.189	1488	1493	1502	rBV	1844351	2536060	26.20%	4.750%
28	10.299	1502	1511	1517	rVV	6865570	9680170	100.00%	18.130%
29	10.360	1517	1521	1529	rVB	269002	384201	3.97%	0.720%
30	10.585	1551	1558	1561	rBV	69967	109770	1.13%	0.206%
31	10.640	1561	1567	1576	rVV	3439251	4348300	44.92%	8.144%
32	10.780	1582	1590	1593	rBV2	98379	160503	1.66%	0.301%
33	10.854	1593	1602	1606	rBV3	85668	148890	1.54%	0.279%
34	10.957	1613	1619	1623	rBV	155898	220519	2.28%	0.413%
35	11.079	1634	1639	1642	rBV	298397	414127	4.28%	0.776%
36	11.189	1650	1657	1660	rBV	99821	143558	1.48%	0.269%

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Instrument :
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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
 Title : SW846 8260

37	11.299	1671	1675	1683	rBV	426159	567144	5.86%	1.062%
38	11.378	1683	1688	1695	rVV	1666037	2893378	29.89%	5.419%
39	11.451	1695	1700	1702	rVV	751345	1019075	10.53%	1.909%
40	11.475	1702	1704	1710	rVB	380757	399955	4.13%	0.749%
41	11.610	1721	1726	1733	rVB	721823	925866	9.56%	1.734%
42	11.713	1740	1743	1745	rBV	90369	103748	1.07%	0.194%
43	11.750	1745	1749	1759	rVB	3043906	3639859	37.60%	6.817%
44	11.890	1769	1772	1776	rVB	78279	99648	1.03%	0.187%
45	11.969	1782	1785	1788	rVV	95932	112325	1.16%	0.210%
46	12.018	1788	1793	1799	rVV3	271086	428919	4.43%	0.803%
47	12.085	1799	1804	1811	rVV	803659	1097966	11.34%	2.056%
48	12.170	1815	1818	1825	rVB2	86580	123565	1.28%	0.231%
49	12.244	1825	1830	1832	rBV2	509052	707869	7.31%	1.326%
50	12.268	1832	1834	1837	rVV	265534	269225	2.78%	0.504%
51	12.317	1837	1842	1848	rVB3	424776	687744	7.10%	1.288%
52	12.420	1848	1859	1863	rBV	424229	571399	5.90%	1.070%
53	12.463	1863	1866	1872	rVB	150313	210964	2.18%	0.395%
54	12.530	1872	1877	1879	rBV	226979	309268	3.19%	0.579%
55	12.555	1879	1881	1884	rVV	205706	221073	2.28%	0.414%
56	12.609	1886	1890	1893	rVV	303251	358887	3.71%	0.672%
57	12.695	1899	1904	1913	rBV3	171833	413011	4.27%	0.774%
58	12.817	1917	1924	1926	rBV4	73148	115622	1.19%	0.217%
59	12.920	1937	1941	1944	rVV	165850	233704	2.41%	0.438%
60	12.963	1944	1948	1954	rVB	285360	471059	4.87%	0.882%
61	13.042	1954	1961	1964	rBV4	116504	199520	2.06%	0.374%
62	13.164	1974	1981	1985	rBV2	222814	324111	3.35%	0.607%
63	13.262	1992	1997	1999	rBV	364831	472257	4.88%	0.884%
64	13.292	1999	2002	2007	rVB2	432902	585382	6.05%	1.096%
65	13.384	2012	2017	2019	rVV2	89376	122879	1.27%	0.230%
66	13.420	2019	2023	2032	rVB	235313	406848	4.20%	0.762%
67	13.499	2032	2036	2039	rBV2	65556	100043	1.03%	0.187%
68	13.585	2046	2050	2055	rBV3	93789	164546	1.70%	0.308%
69	13.664	2060	2063	2067	rBV3	64882	103625	1.07%	0.194%
70	13.774	2077	2081	2087	rVB	193094	277484	2.87%	0.520%
71	13.847	2087	2093	2099	rVB6	146965	289465	2.99%	0.542%
72	13.926	2099	2106	2111	rBV6	103179	209327	2.16%	0.392%
73	14.036	2120	2124	2127	rBV	351233	343722	3.55%	0.644%
74	14.225	2150	2155	2161	rVB2	208870	354239	3.66%	0.663%
75	14.475	2192	2196	2206	rVV5	131233	297089	3.07%	0.556%
76	14.609	2215	2218	2221	rVV	182310	261396	2.70%	0.490%

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ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

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\82X040225W.M
Title : SW846 8260

77	14.755	2238	2242	2245	rBV	327036	357096	3.69%	0.669%
78	15.182	2307	2312	2314	rBV2	78514	118614	1.23%	0.222%
79	15.213	2314	2317	2320	rVV	120032	160227	1.66%	0.300%
80	15.487	2357	2362	2368	rBV	264425	381439	3.94%	0.714%
81	16.371	2502	2507	2514	rVB	122556	203011	2.10%	0.380%

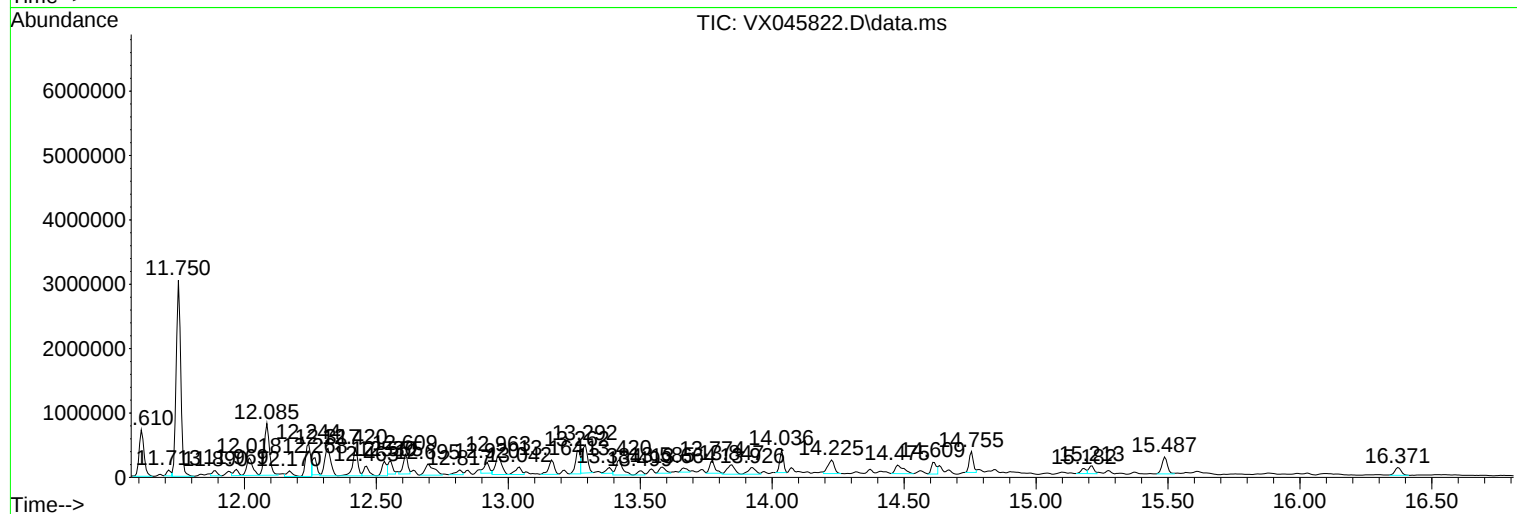
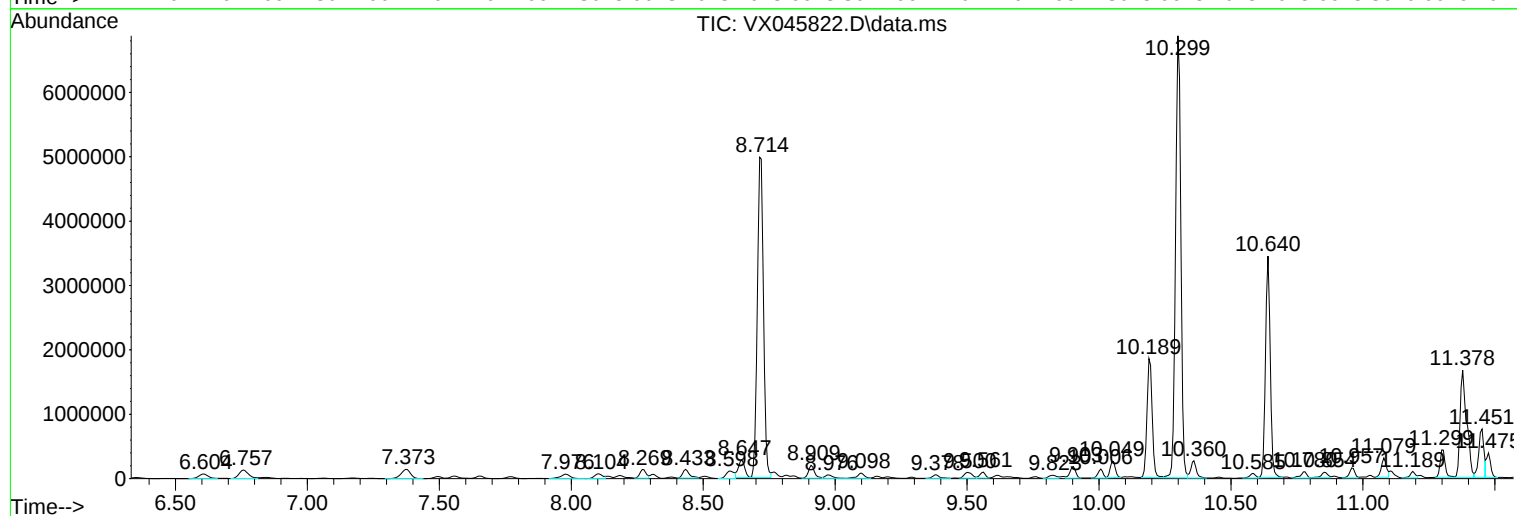
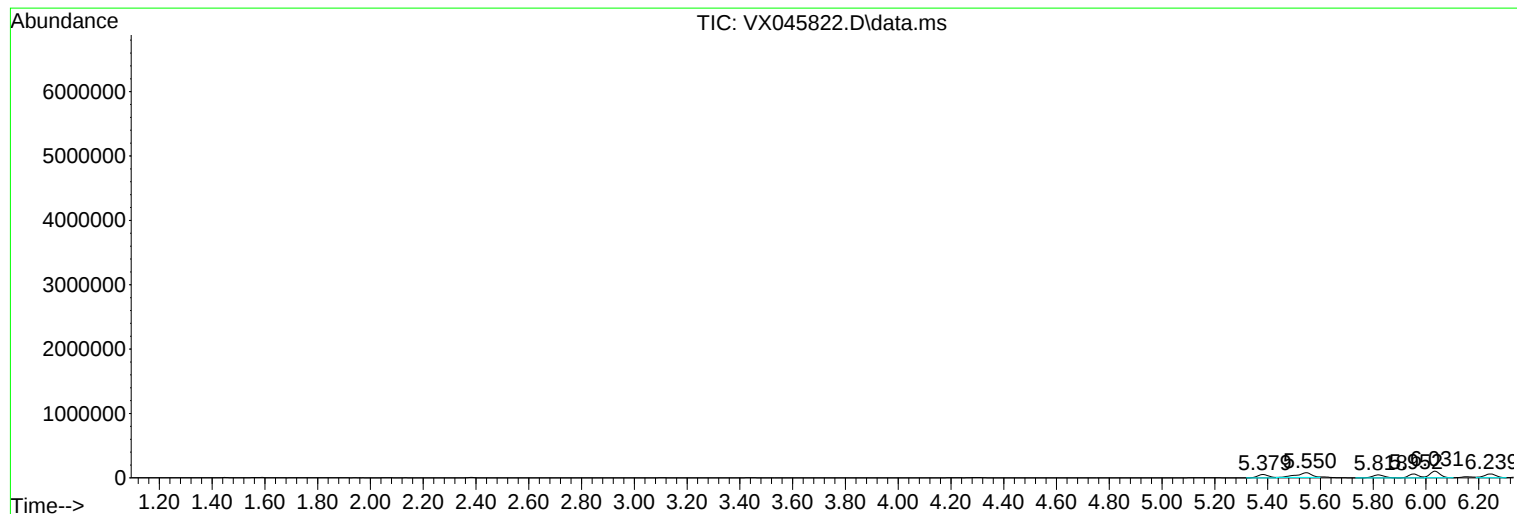
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Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

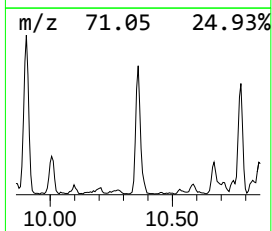
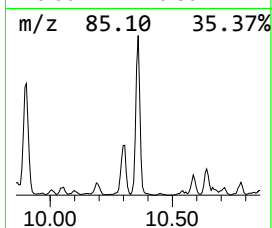
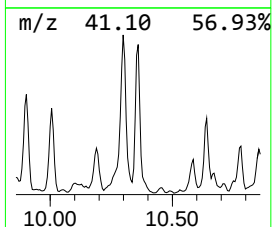
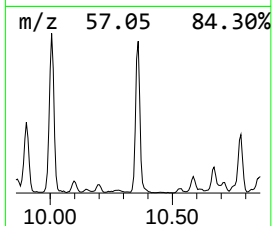
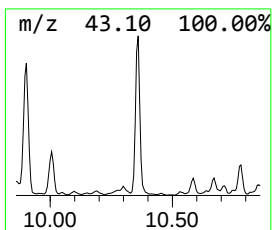
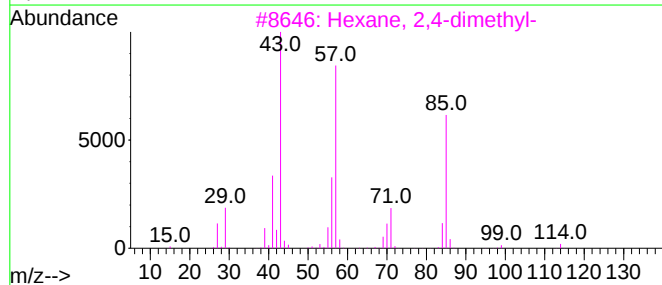
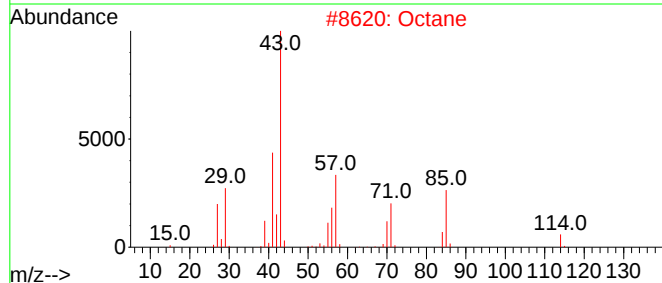
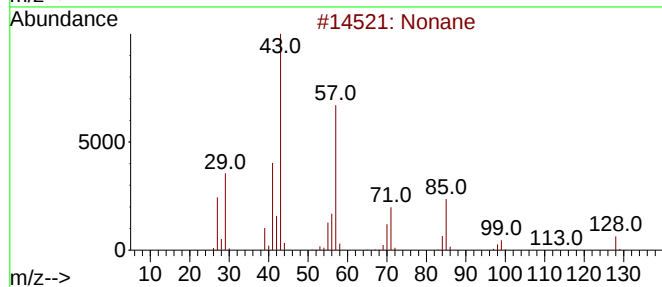
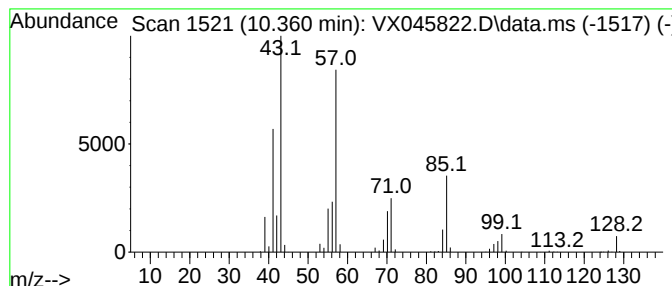
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
Quant Title : SW846 8260

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

Peak Number 1 Nonane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	55.72 ug/l	384201	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	94
2			Octane	114	C8H18	000111-65-9	59
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	53
4			Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	53
5			2-Nonanol, trimethylacetate	228	C14H28O2	1000463-21-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VO41725\
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Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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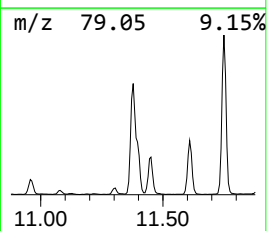
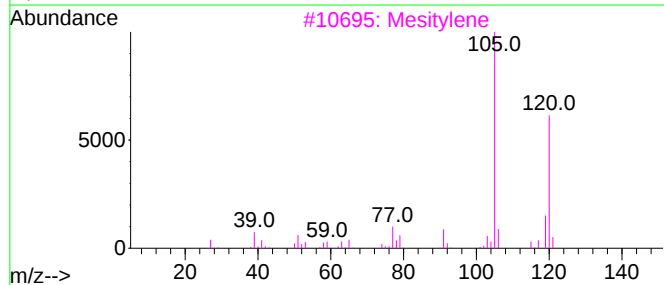
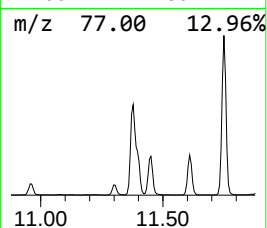
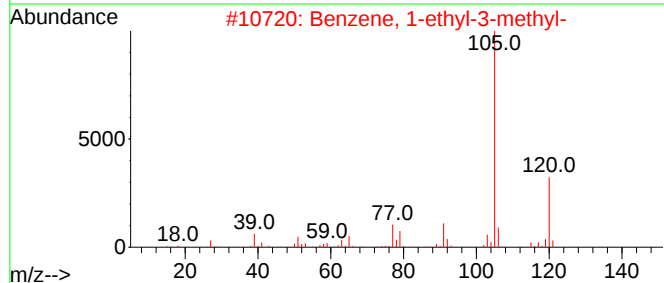
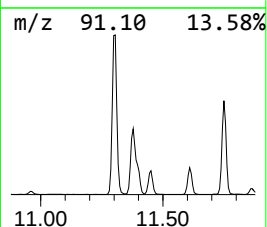
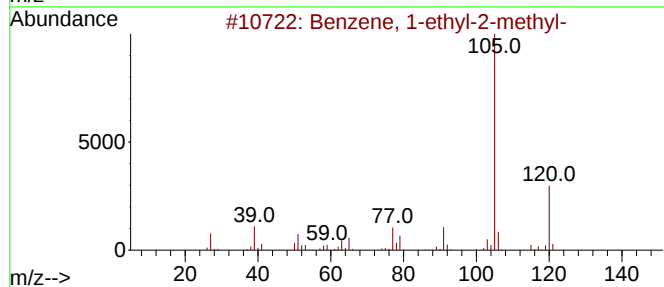
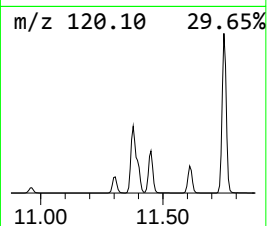
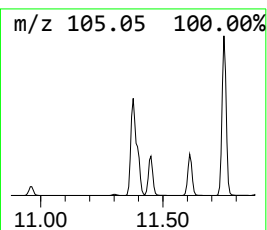
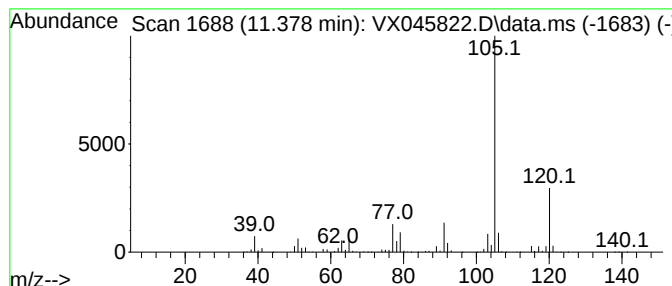
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	337.29 ug/l	2893380	1,4-Dichlorobenzene-d4	12.018

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	94
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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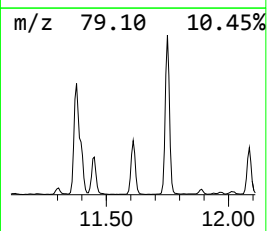
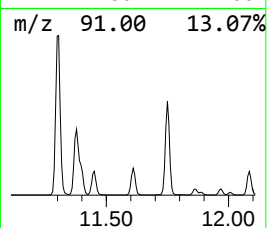
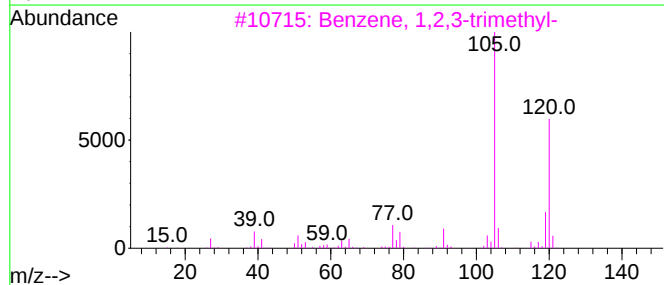
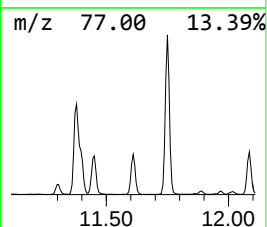
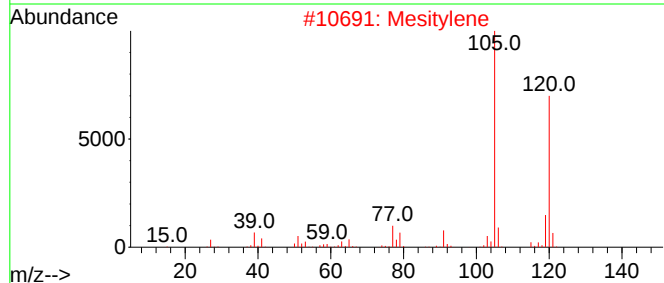
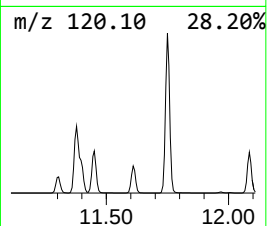
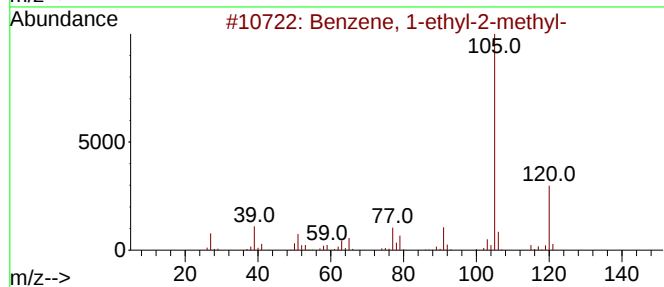
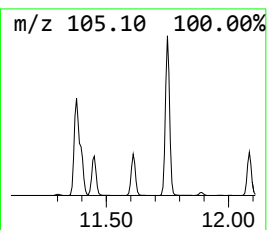
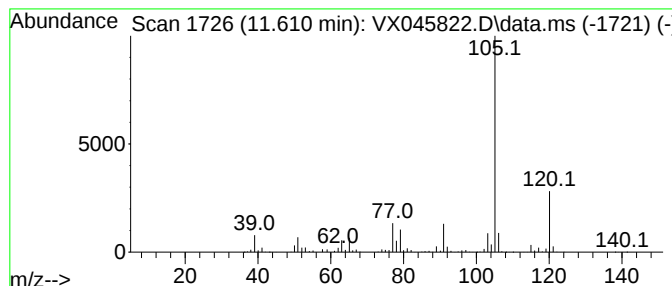
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	107.93 ug/l	925866	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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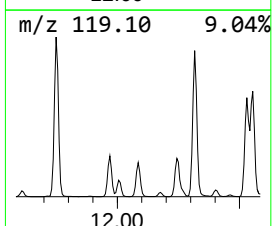
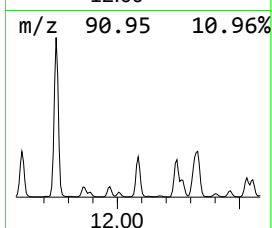
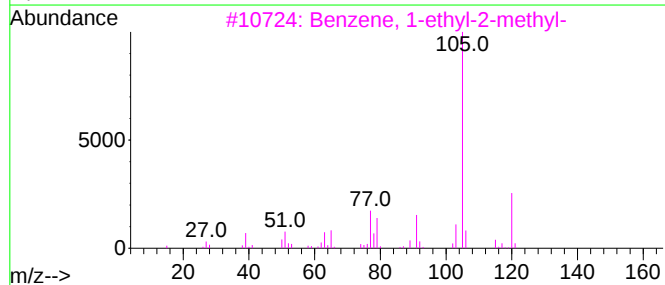
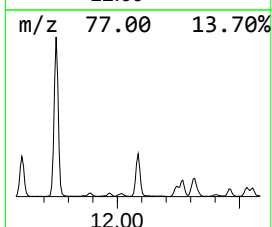
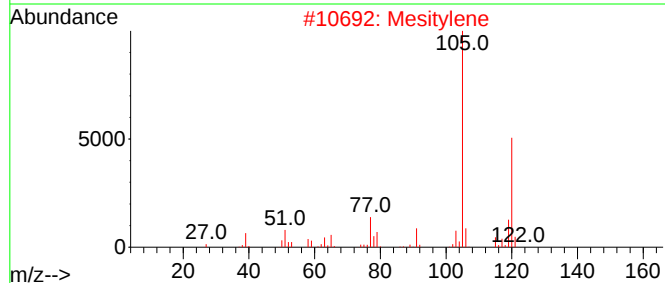
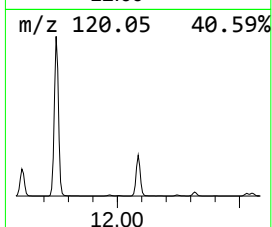
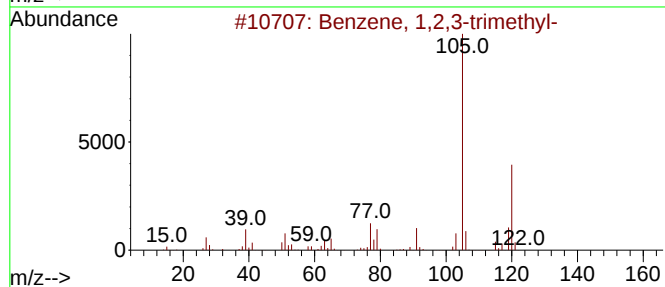
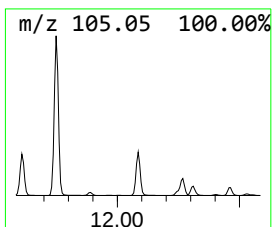
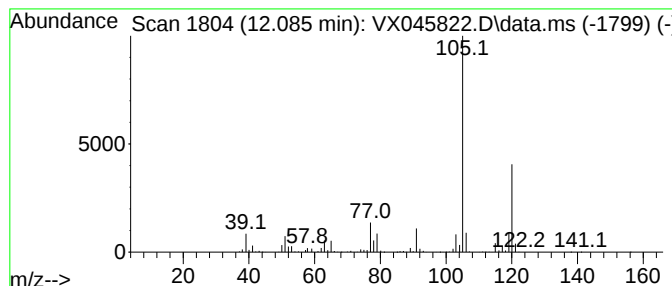
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	127.99 ug/l	1097970	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VO41725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

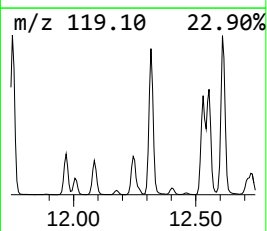
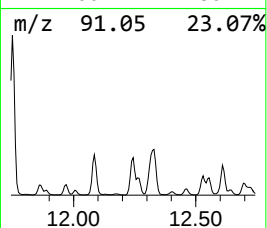
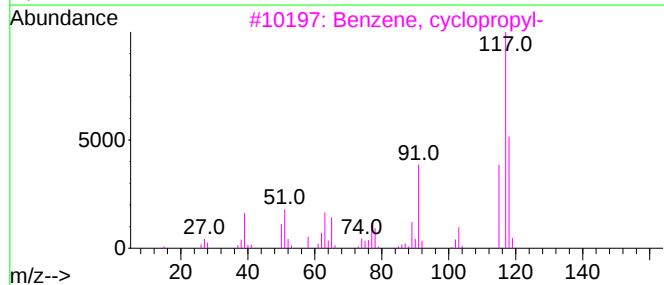
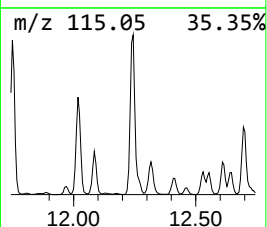
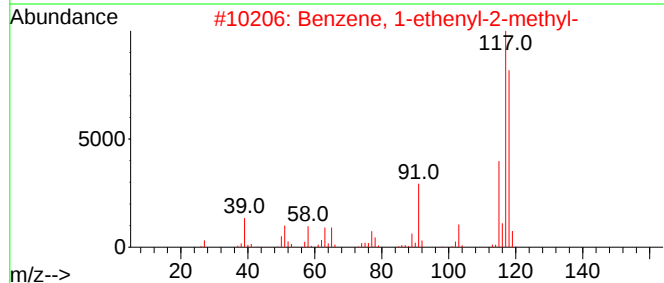
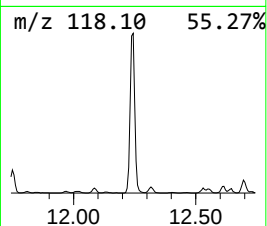
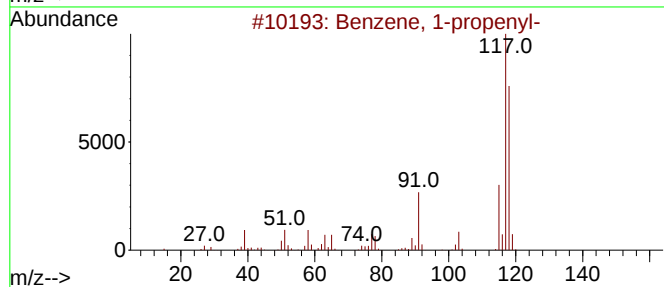
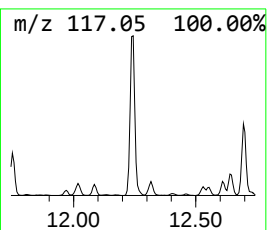
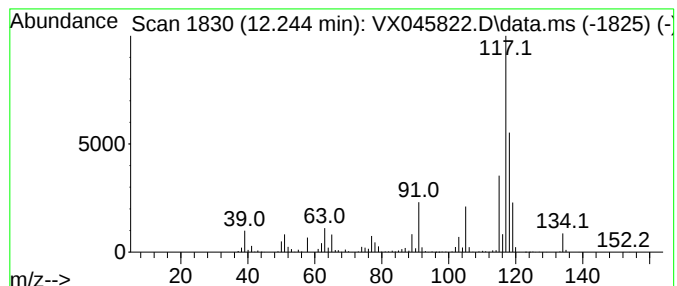
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-propenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	82.52 ug/l	707869	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4			Indane	118	C9H10	000496-11-7	70
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

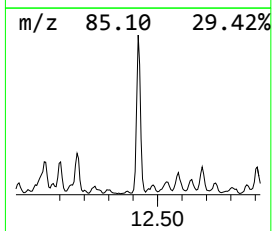
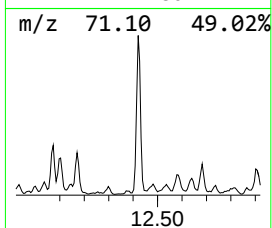
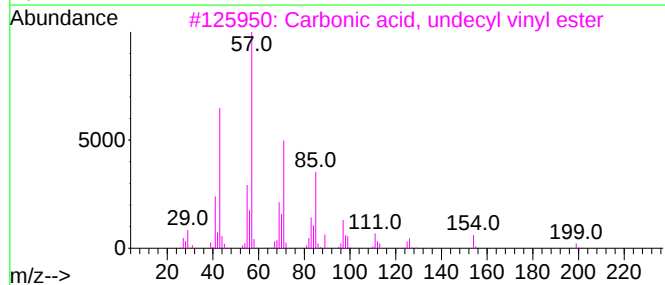
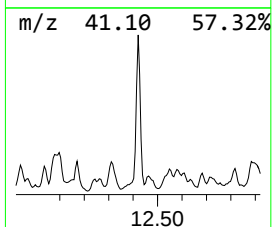
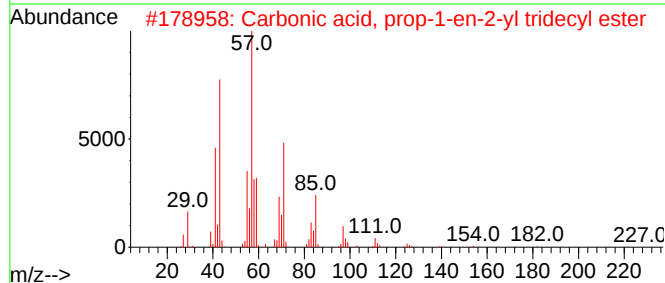
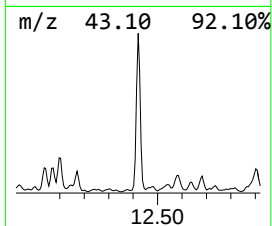
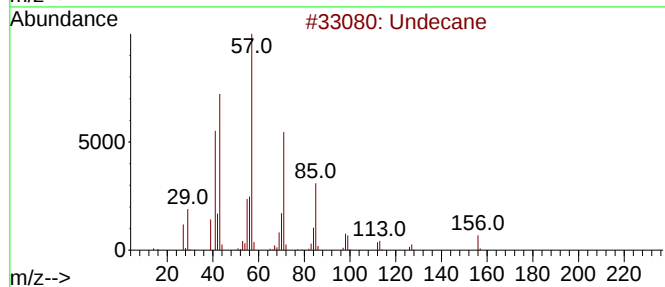
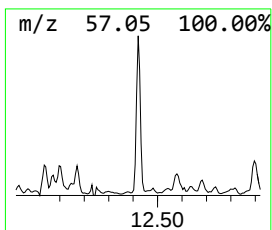
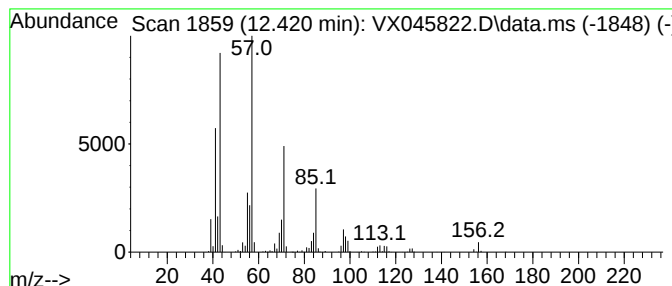
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Quant Title : SW846 8260

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

Peak Number 6 Undecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.420	66.61 ug/l	571399	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	93
2			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	90
3			Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	83
4			Decane	142	C10H22	000124-18-5	78
5			Heptadecane	240	C17H36	000629-78-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

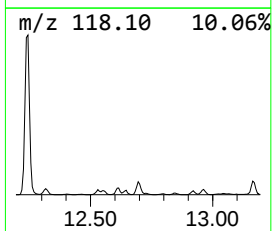
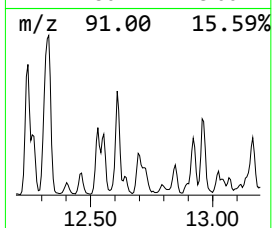
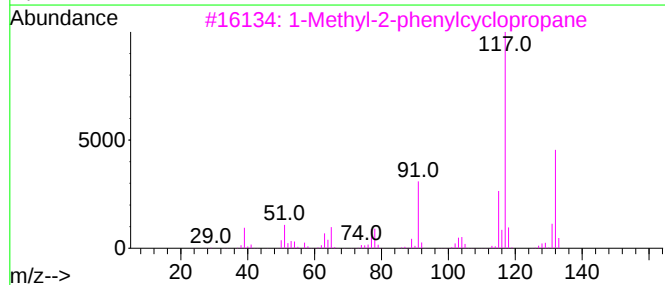
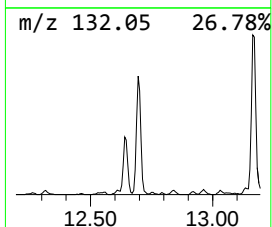
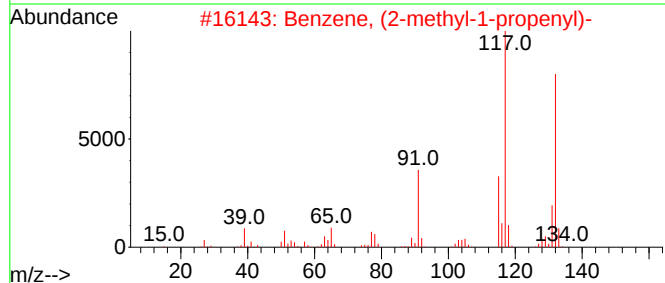
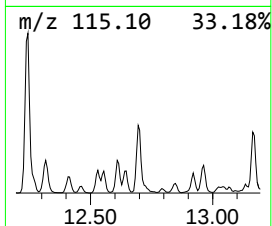
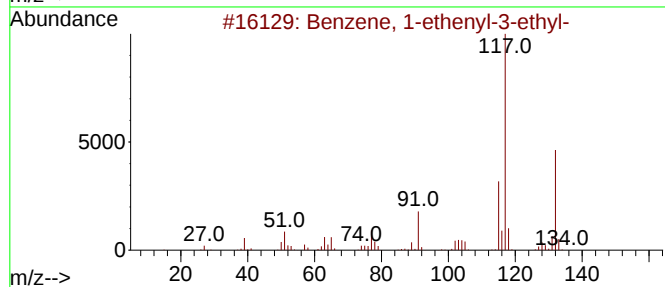
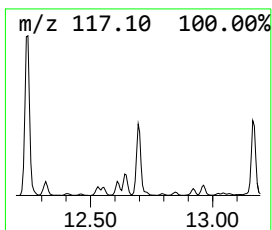
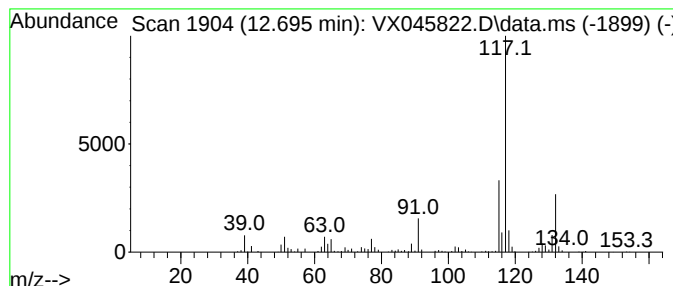
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	48.15 ug/l	413011	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Indan, 1-methyl-	132	C10H12	000767-58-8	87
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

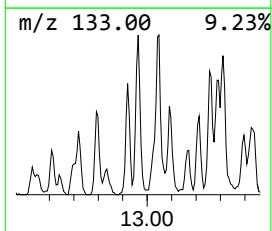
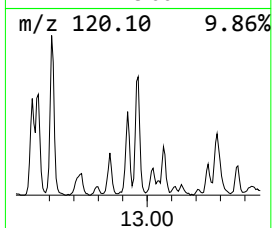
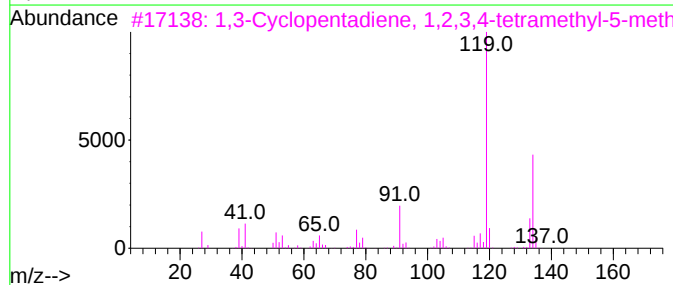
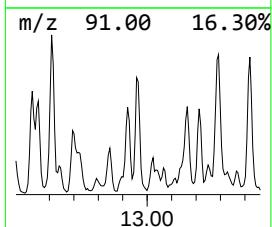
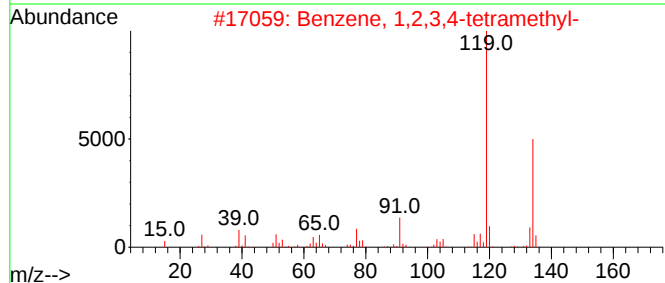
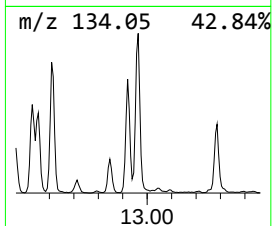
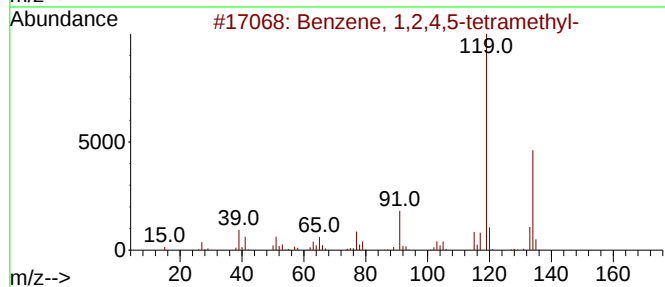
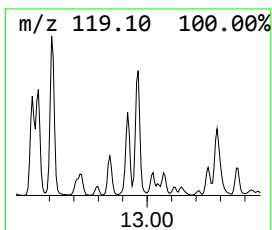
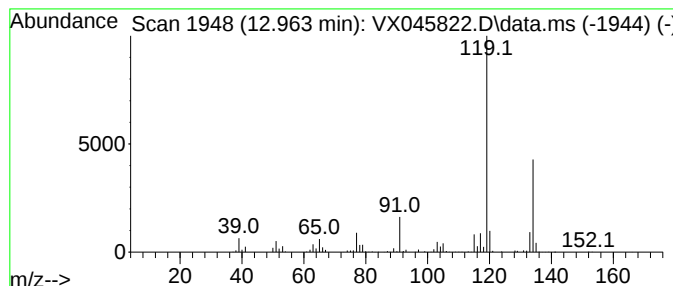
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	54.91 ug/l	471059	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

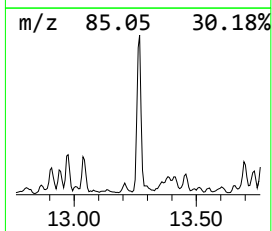
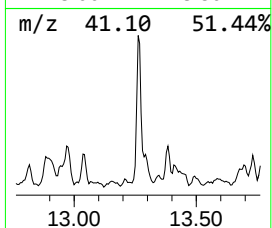
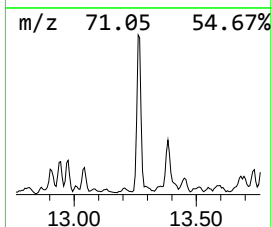
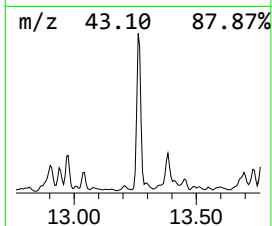
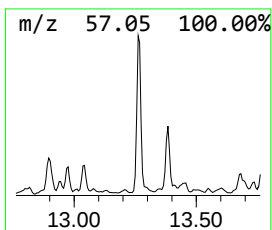
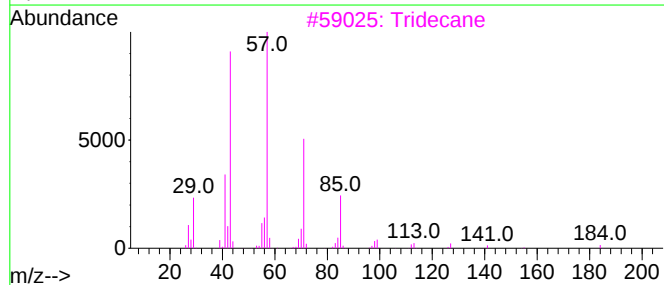
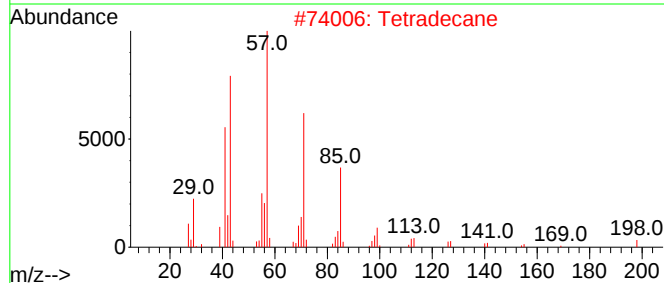
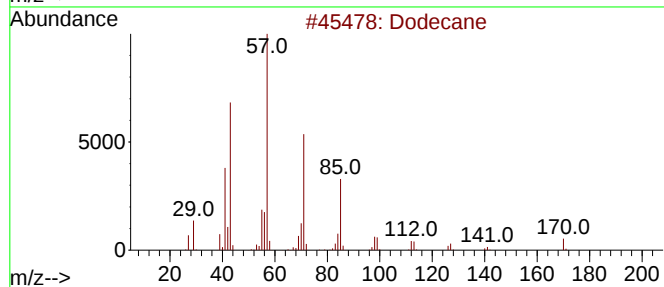
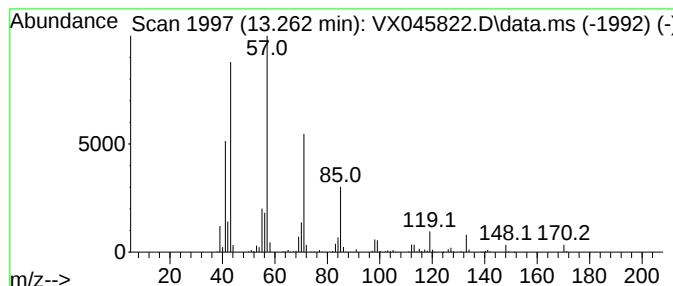
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Quant Title : SW846 8260

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

Peak Number 9 Dodecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.262	55.05 ug/l	472257	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	94
2			Tetradecane	198	C14H30	000629-59-4	80
3			Tridecane	184	C13H28	000629-50-5	72
4			Hexadecane	226	C16H34	000544-76-3	72
5			Undecane	156	C11H24	001120-21-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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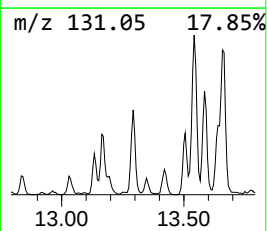
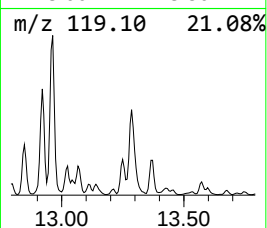
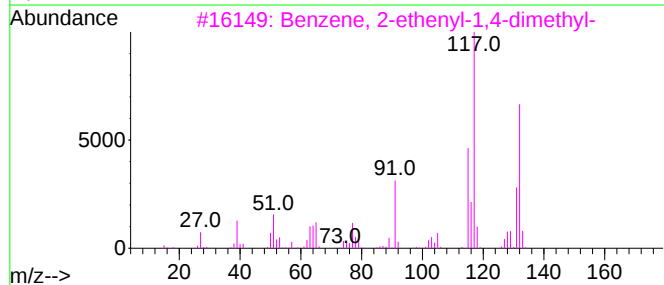
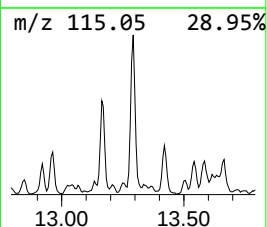
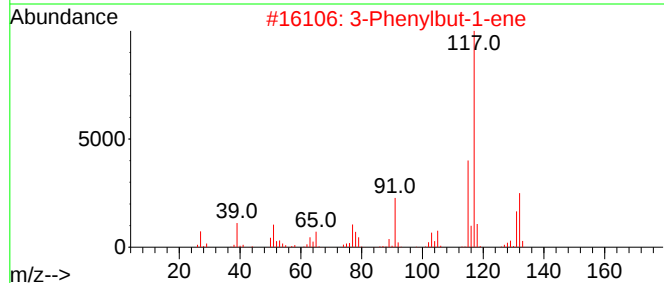
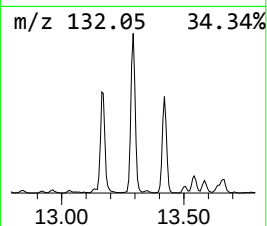
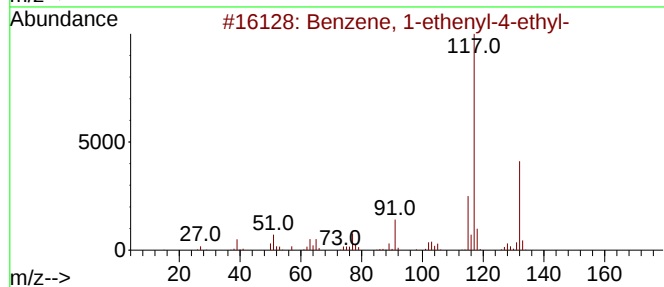
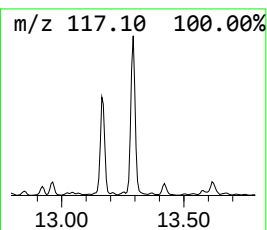
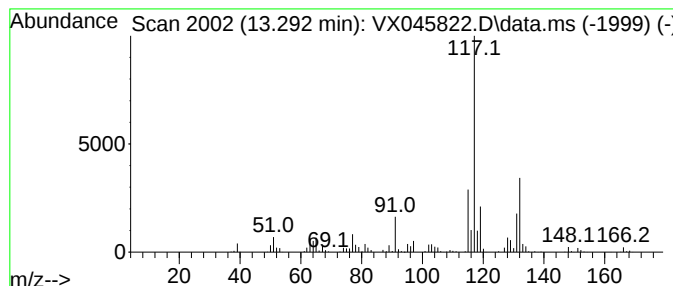
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	68.24 ug/l	585382	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
5			Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX041725\
Data File : VX045822.D
Acq On : 17 Apr 2025 12:49
Operator : JC/MD
Sample : Q1824-04 50X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
285

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.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
Nonane	10.360	55.7	ug/l	384201	3	10.055	344789	50.0
Benzene, 1-ethy...	11.378	337.3	ug/l	2893380	4	12.018	428919	50.0
Benzene, 1-ethy...	11.610	107.9	ug/l	925866	4	12.018	428919	50.0
Benzene, 1,2,3-...	12.085	128.0	ug/l	1097970	4	12.018	428919	50.0
Benzene, 1-prop...	12.244	82.5	ug/l	707869	4	12.018	428919	50.0
Undecane	12.420	66.6	ug/l	571399	4	12.018	428919	50.0
Benzene, 1-ethe...	12.695	48.1	ug/l	413011	4	12.018	428919	50.0
Benzene, 1,2,4,...	12.963	54.9	ug/l	471059	4	12.018	428919	50.0
Dodecane	13.262	55.0	ug/l	472257	4	12.018	428919	50.0
Benzene, 1-ethe...	13.292	68.2	ug/l	585382	4	12.018	428919	50.0