

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
 Data File : VX046601.D
 Acq On : 10 Jun 2025 14:38
 Operator : JC/MD
 Sample : Q2230-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GW-MW901-060425

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\82X060625W.M
 Title : SW846 8260

Signal : TIC: VX046601.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.392	206	214	229	rBV	29383	59116	7.34%	0.544%
2	4.580	567	573	583	rBV2	3828	12012	1.49%	0.111%
3	5.105	650	659	670	rBV2	5901	18955	2.35%	0.174%
4	5.397	696	707	725	rBV	88342	280373	34.81%	2.581%
5	5.562	725	734	759	rVB2	120495	367407	45.61%	3.382%
6	5.970	791	801	813	rBV	96150	267617	33.22%	2.463%
7	6.769	923	932	952	rBV	218112	550108	68.29%	5.063%
8	7.385	1025	1033	1041	rVB4	5006	12634	1.57%	0.116%
9	8.653	1234	1241	1249	rBV	497402	805537	100.00%	7.414%
10	8.726	1249	1253	1260	rVB	15551	26644	3.31%	0.245%
11	10.055	1465	1471	1480	rBV	540193	743604	92.31%	6.844%
12	10.195	1488	1494	1502	rBB	62115	82038	10.18%	0.755%
13	10.299	1502	1511	1520	rBV	239107	334456	41.52%	3.078%
14	10.640	1562	1567	1580	rVB	323945	406944	50.52%	3.746%
15	10.963	1613	1620	1627	rBV	21545	28450	3.53%	0.262%
16	11.079	1634	1639	1654	rBV	480360	601064	74.62%	5.532%
17	11.305	1671	1676	1683	rBV	42578	53767	6.67%	0.495%
18	11.378	1683	1688	1696	rVV	172394	345847	42.93%	3.183%
19	11.451	1696	1700	1709	rVB	120315	152691	18.96%	1.405%
20	11.610	1721	1726	1740	rVB	191830	239196	29.69%	2.202%
21	11.750	1744	1749	1763	rVB	440898	545942	67.77%	5.025%
22	11.890	1768	1772	1778	rVB	17085	22624	2.81%	0.208%
23	11.969	1778	1785	1788	rBV	33829	40964	5.09%	0.377%
24	12.018	1788	1793	1799	rVV	611675	759761	94.32%	6.993%
25	12.085	1799	1804	1816	rVB	228770	292222	36.28%	2.690%
26	12.244	1824	1830	1832	rBV2	146582	200586	24.90%	1.846%
27	12.268	1832	1834	1837	rVV	62916	64969	8.07%	0.598%
28	12.317	1837	1842	1851	rVB3	97184	157984	19.61%	1.454%
29	12.402	1851	1856	1861	rVB	15891	21684	2.69%	0.200%
30	12.463	1861	1866	1871	rBV	49988	61622	7.65%	0.567%
31	12.530	1872	1877	1879	rBV	61608	85479	10.61%	0.787%
32	12.555	1879	1881	1885	rVB	59040	61838	7.68%	0.569%
33	12.610	1885	1890	1893	rBV	89570	109919	13.65%	1.012%
34	12.640	1893	1895	1900	rVB	29023	35102	4.36%	0.323%
35	12.695	1900	1904	1917	rVB2	118698	195774	24.30%	1.802%
36	12.792	1917	1920	1924	rBV2	10233	13177	1.64%	0.121%

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Integrator: RTE

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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

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37	12.847	1924	1929	1933	rVB2	41287	55591	6.90%	0.512%
38	12.920	1933	1941	1944	rBV	56093	72415	8.99%	0.667%
39	12.963	1944	1948	1954	rVB	78169	99673	12.37%	0.917%
40	13.024	1954	1958	1960	rBV	19928	28954	3.59%	0.266%
41	13.067	1963	1965	1968	rVB	10005	9540	1.18%	0.088%
42	13.134	1973	1976	1978	rBV2	16842	22283	2.77%	0.205%
43	13.164	1978	1981	1985	rVB	89734	106358	13.20%	0.979%
44	13.207	1985	1988	1992	rVB3	16312	20400	2.53%	0.188%
45	13.250	1992	1995	1997	rBV3	19281	25315	3.14%	0.233%
46	13.292	1997	2002	2007	rVV2	214606	289968	36.00%	2.669%
47	13.365	2012	2014	2018	rVV	13934	17564	2.18%	0.162%
48	13.420	2018	2023	2031	rVB	171615	217564	27.01%	2.002%
49	13.506	2031	2037	2039	rBV2	21846	30011	3.73%	0.276%
50	13.542	2039	2043	2046	rVV	47877	68171	8.46%	0.627%
51	13.585	2046	2050	2054	rVV3	53111	97511	12.11%	0.898%
52	13.640	2055	2059	2060	rVV2	27961	43567	5.41%	0.401%
53	13.658	2060	2062	2069	rVB2	56449	82653	10.26%	0.761%
54	13.774	2074	2081	2088	rVV	49116	73157	9.08%	0.673%
55	13.847	2089	2093	2099	rVB	68668	94831	11.77%	0.873%
56	13.920	2099	2105	2110	rBV2	73015	111199	13.80%	1.023%
57	13.969	2110	2113	2117	rVV	27922	34136	4.24%	0.314%
58	14.000	2117	2118	2125	rVB5	7734	12049	1.50%	0.111%
59	14.073	2125	2130	2134	rBV	41163	50641	6.29%	0.466%
60	14.225	2147	2155	2162	rBV	153163	257845	32.01%	2.373%
61	14.317	2167	2170	2175	rVV	20008	23779	2.95%	0.219%
62	14.371	2175	2179	2183	rVB	42854	53377	6.63%	0.491%
63	14.414	2183	2186	2191	rVB2	9764	14303	1.78%	0.132%
64	14.475	2191	2196	2202	rBV2	98219	138097	17.14%	1.271%
65	14.615	2214	2219	2220	rBV	54588	71674	8.90%	0.660%
66	14.634	2220	2222	2225	rVV	93078	109044	13.54%	1.004%
67	14.670	2225	2228	2233	rVB	35948	47576	5.91%	0.438%
68	14.725	2233	2237	2241	rBV	11370	14385	1.79%	0.132%
69	14.780	2241	2246	2249	rBV	43286	59863	7.43%	0.551%
70	14.816	2249	2252	2254	rVV2	20968	27405	3.40%	0.252%
71	14.841	2254	2256	2261	rVB3	18004	23421	2.91%	0.216%
72	14.902	2261	2266	2273	rBV3	19393	38455	4.77%	0.354%
73	15.048	2283	2290	2294	rBV3	9523	15452	1.92%	0.142%
74	15.182	2305	2312	2314	rBV	33234	51308	6.37%	0.472%
75	15.249	2319	2323	2330	rVB9	5676	12361	1.53%	0.114%
76	15.371	2338	2343	2347	rBV2	12456	19733	2.45%	0.182%

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77	15.475	2353	2360	2364	rBV2	27740	44727	5.55%	0.412%
78	15.524	2364	2368	2373	rVV5	8148	17986	2.23%	0.166%
79	15.609	2377	2382	2386	rVV2	31386	54104	6.72%	0.498%
80	15.646	2386	2388	2397	rVB3	18208	27308	3.39%	0.251%
81	15.841	2412	2420	2425	rBV5	6582	12200	1.51%	0.112%
82	16.085	2455	2460	2470	rBV9	3605	8674	1.08%	0.080%

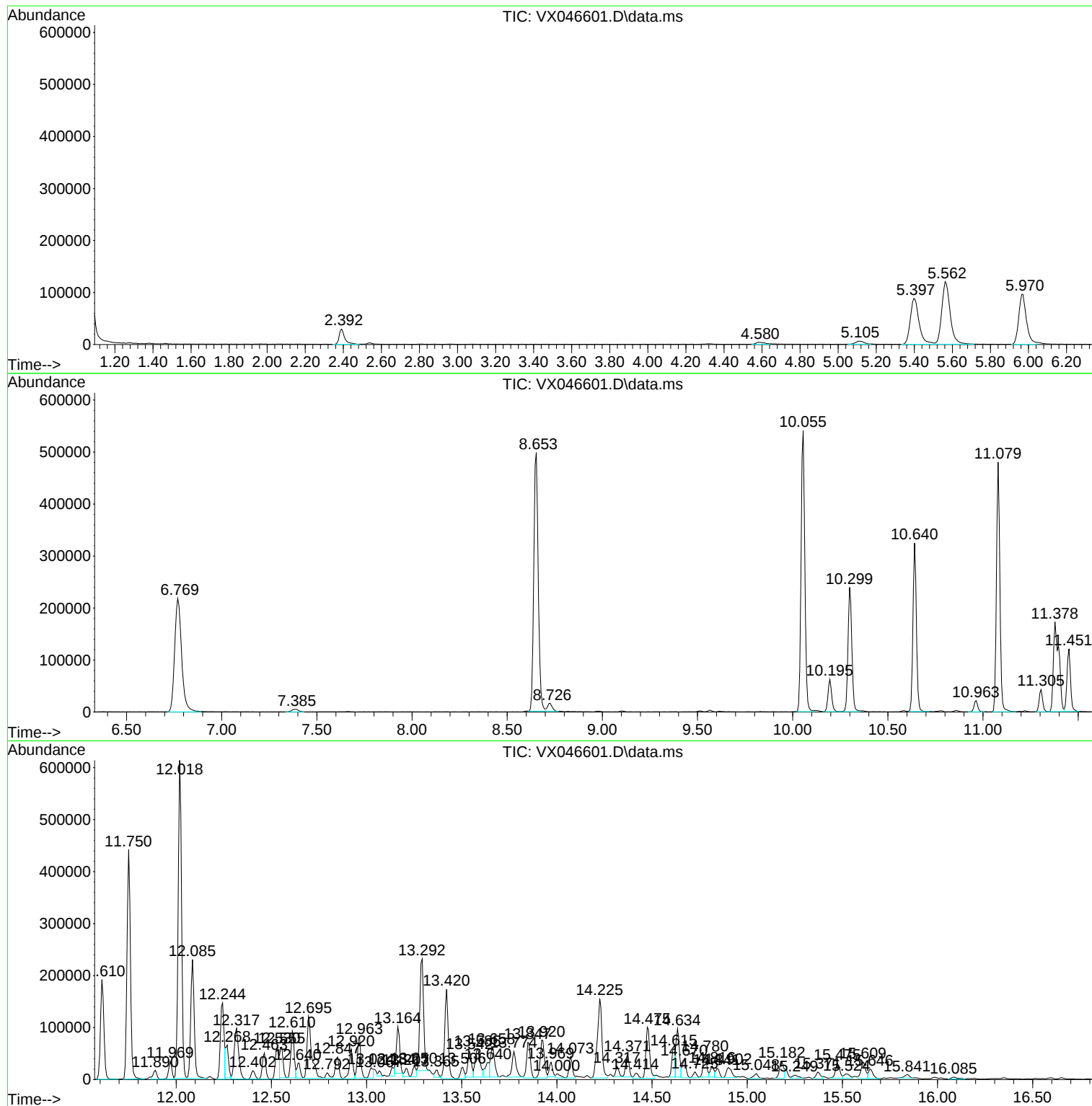
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Quant Title : SW846 8260

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



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Instrument :
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GW-MW901-060425

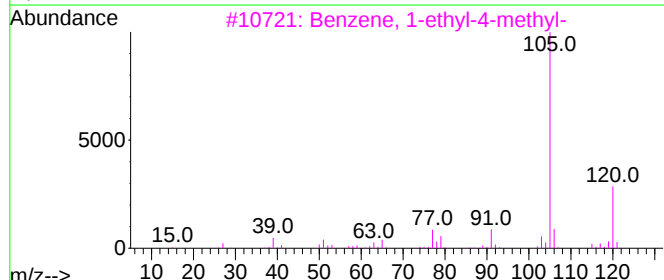
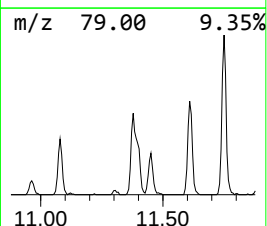
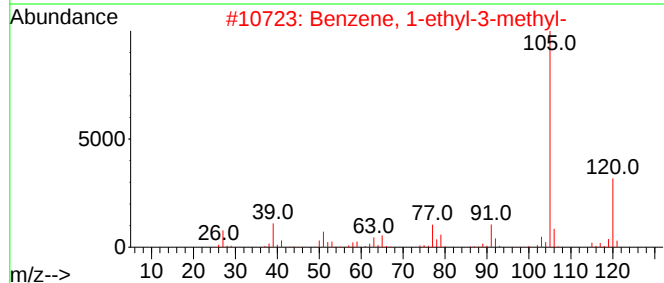
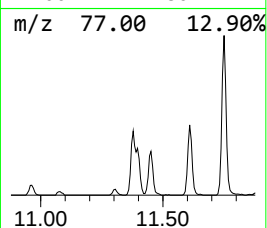
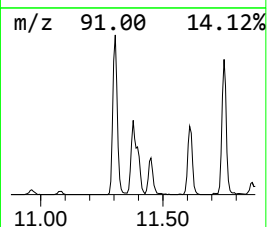
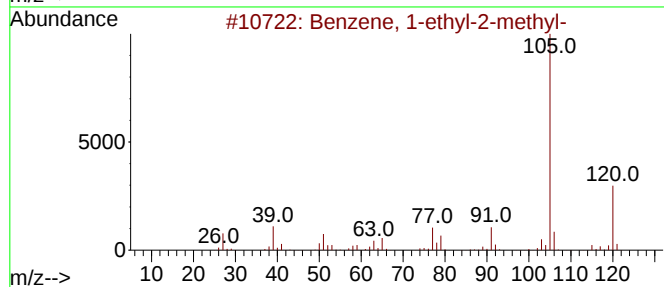
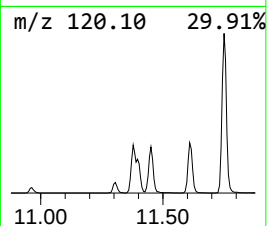
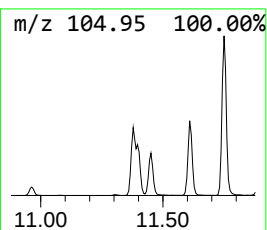
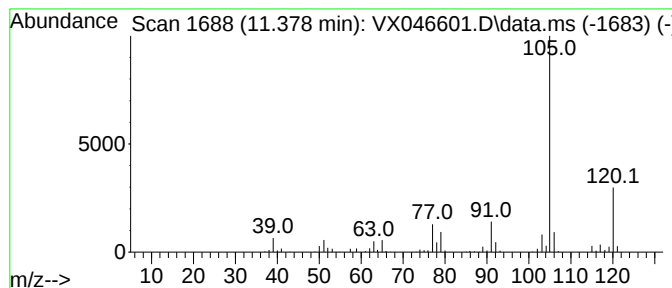
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	22.76 ug/l	345847	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			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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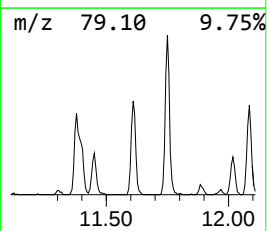
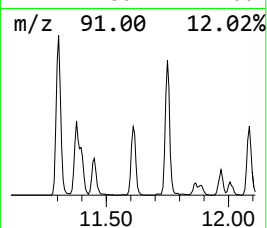
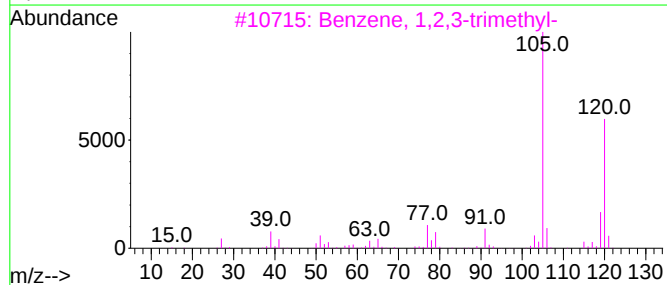
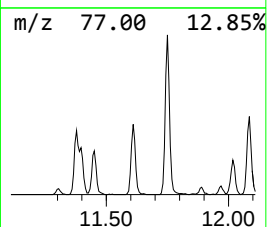
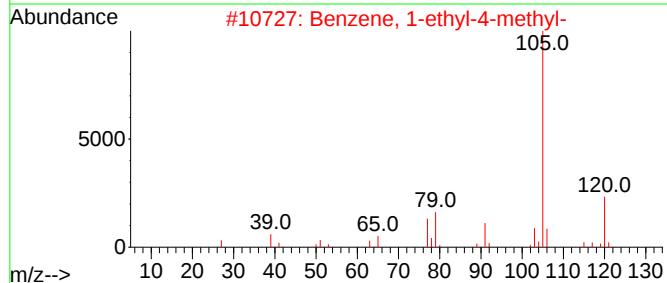
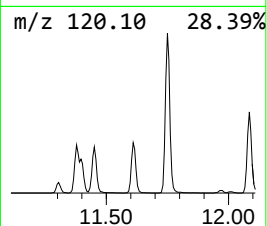
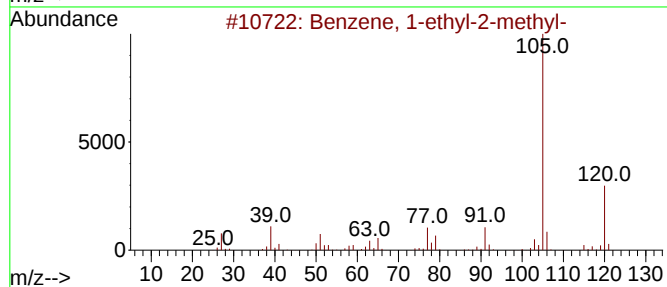
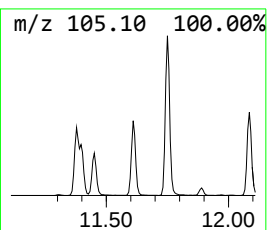
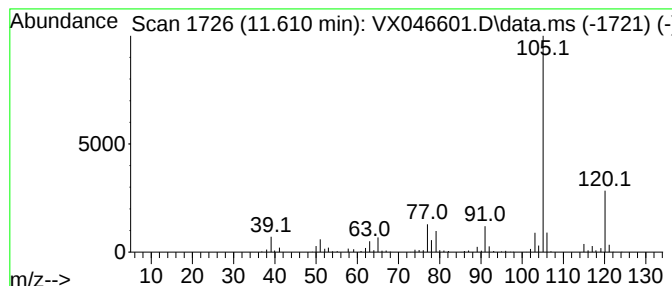
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	15.74 ug/l	239196	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-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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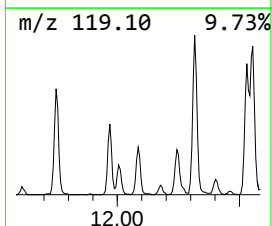
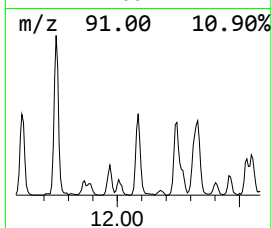
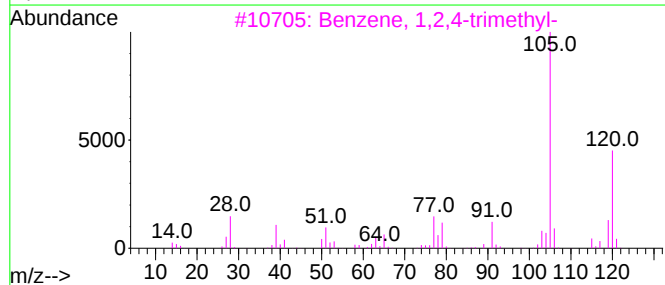
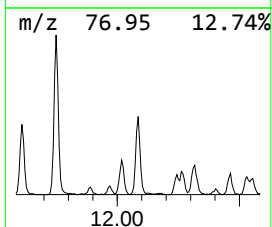
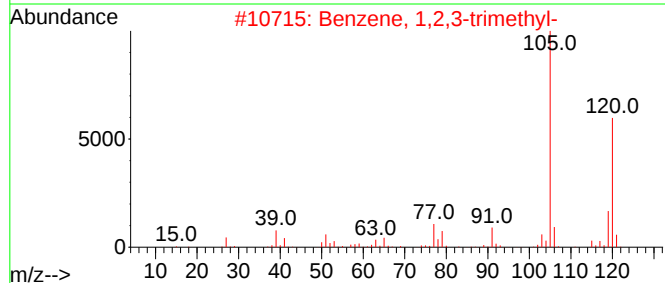
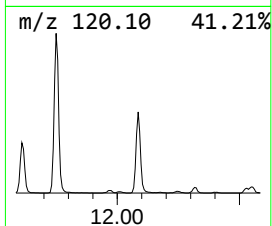
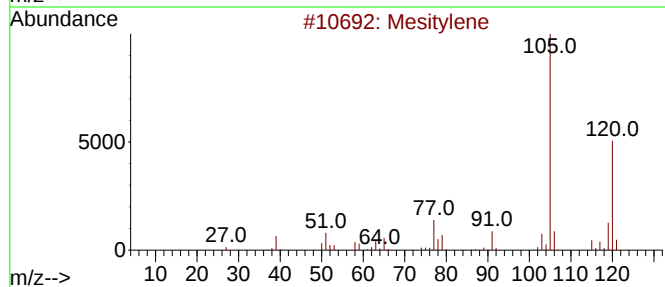
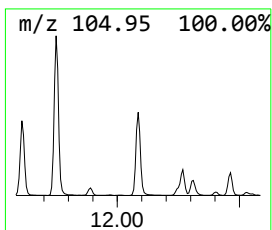
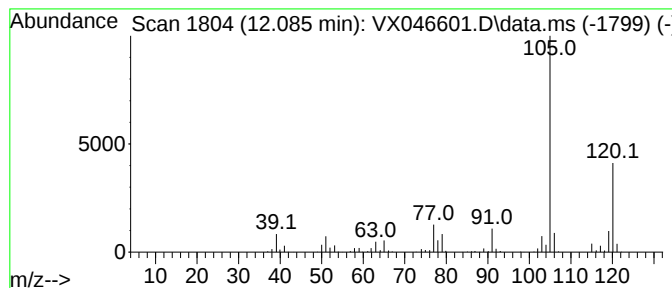
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	19.23 ug/l	292222	1,4-Dichlorobenzene-d4	12.018

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	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



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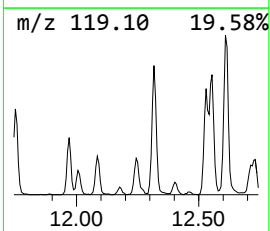
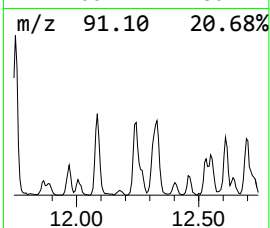
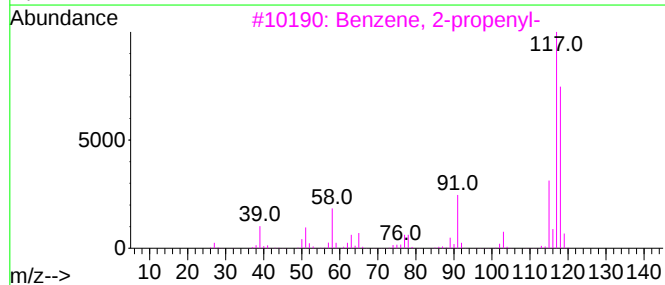
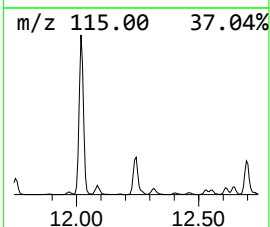
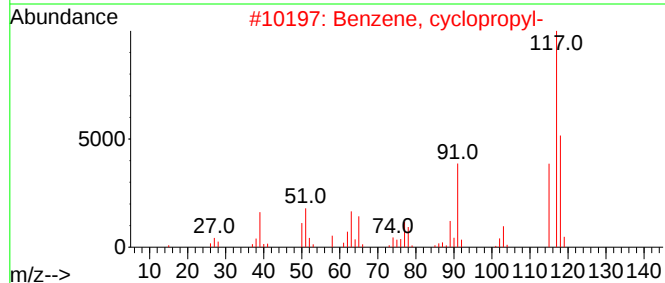
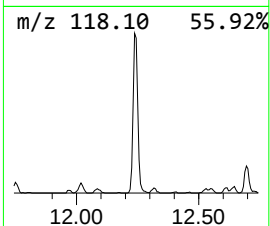
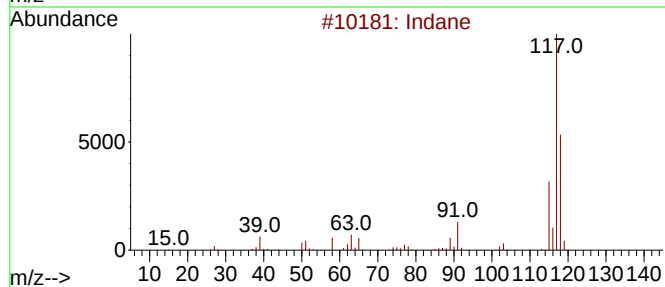
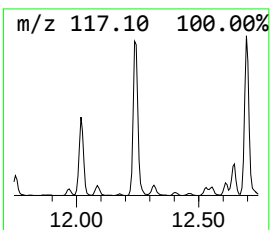
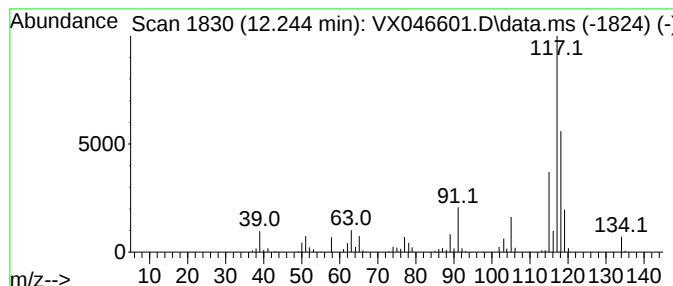
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	13.20 ug/l	200586	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	64
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

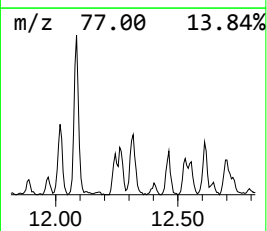
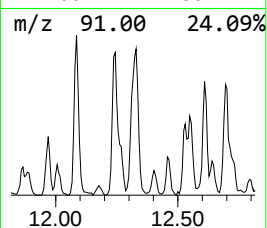
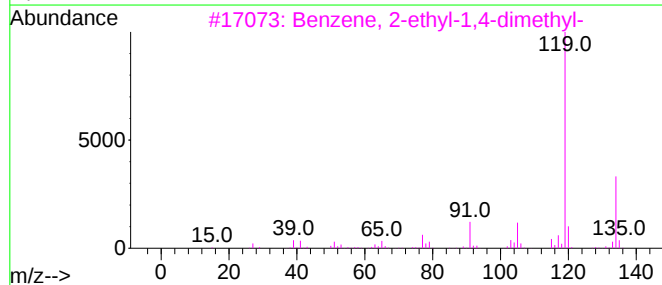
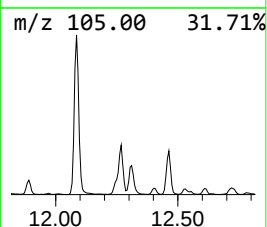
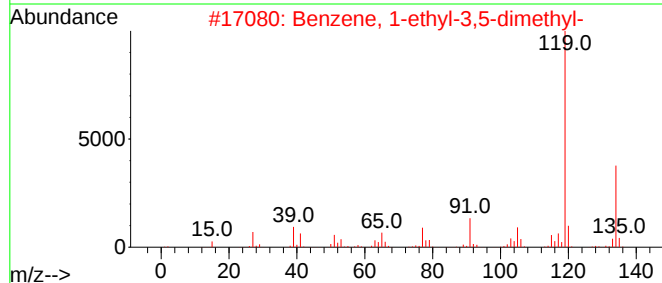
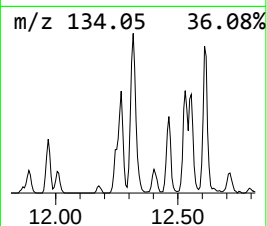
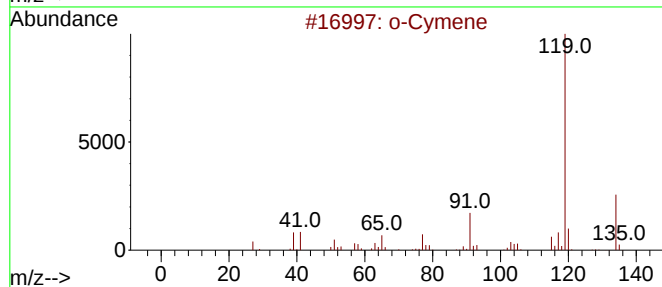
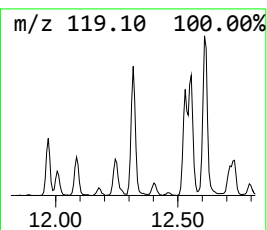
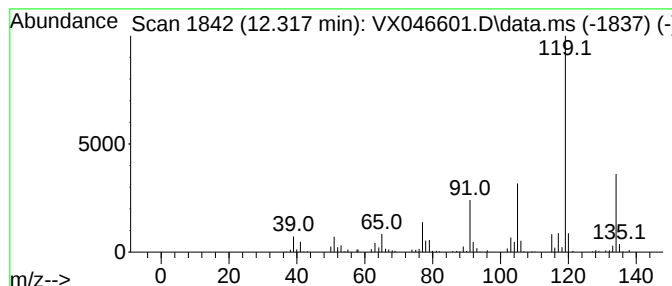
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	10.40 ug/l	157984	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5			p-Cymene	134	C10H14	000099-87-6	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

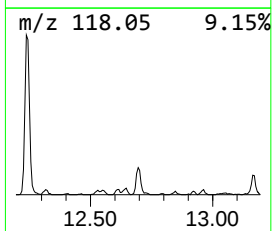
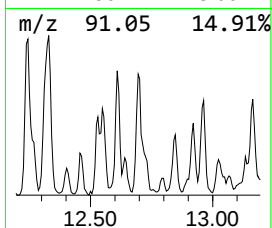
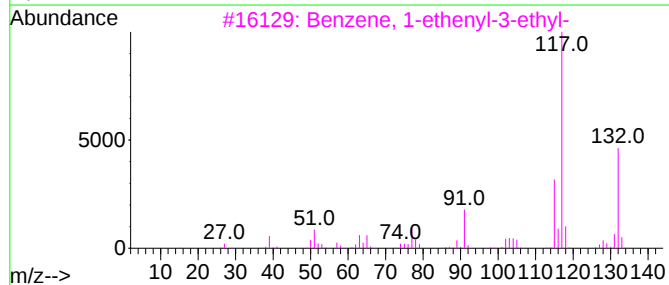
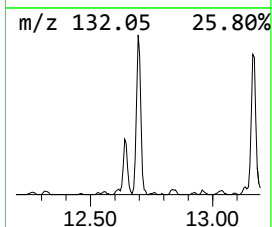
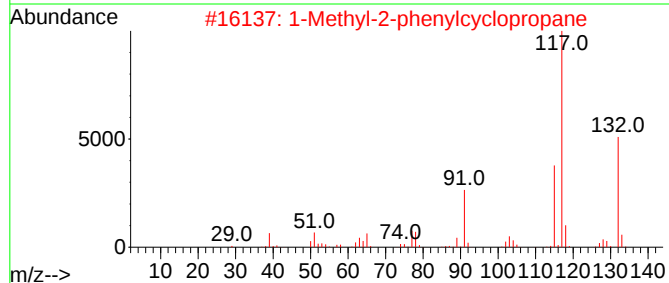
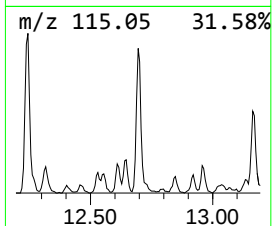
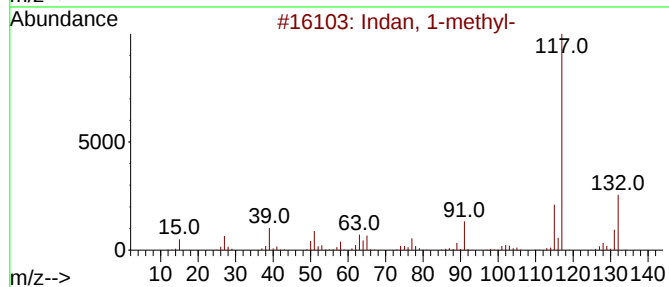
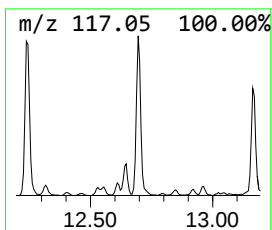
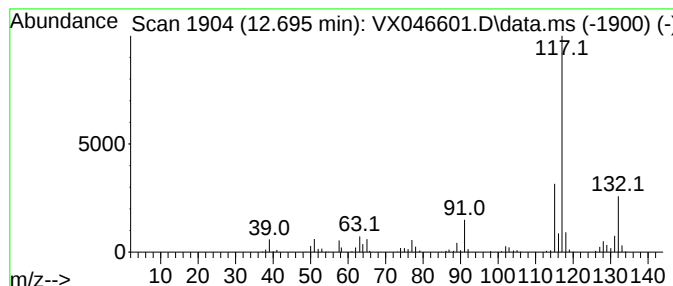
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	12.88 ug/l	195774	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

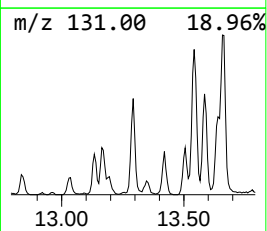
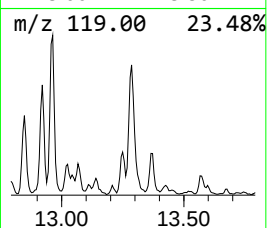
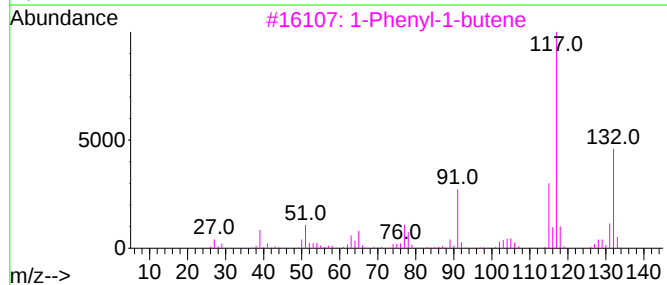
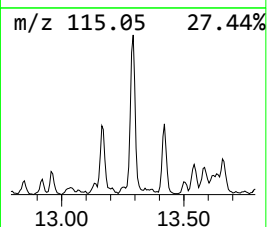
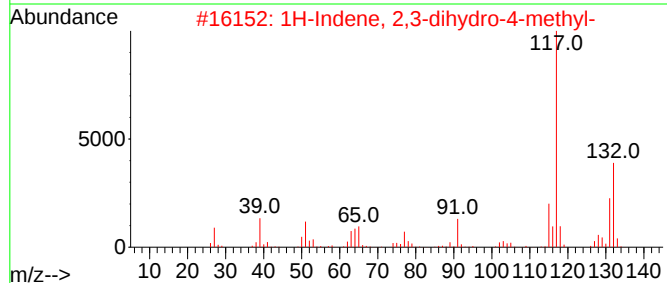
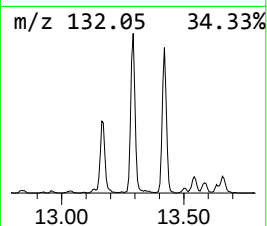
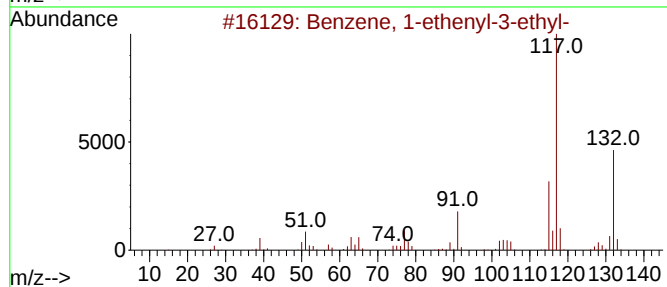
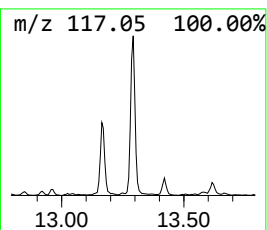
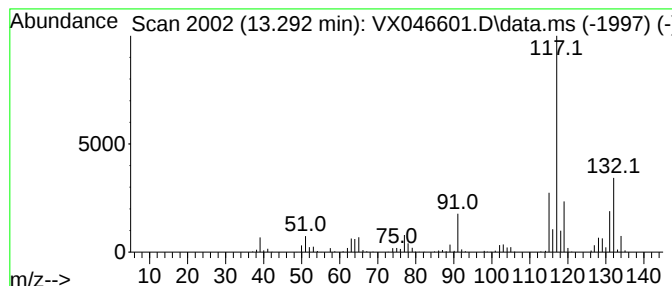
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	19.08 ug/l	289968	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	81
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	81
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	81
4			Indan, 1-methyl-	132	C10H12	000767-58-8	76
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

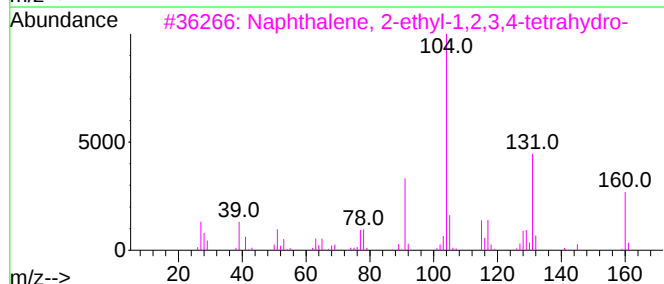
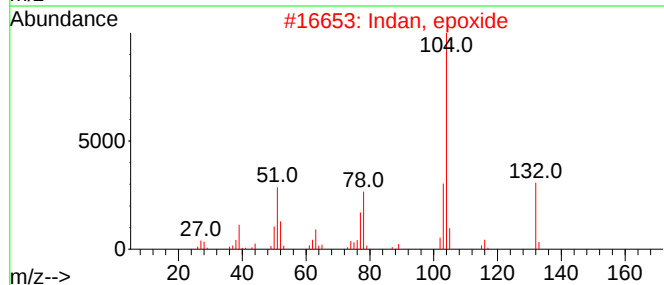
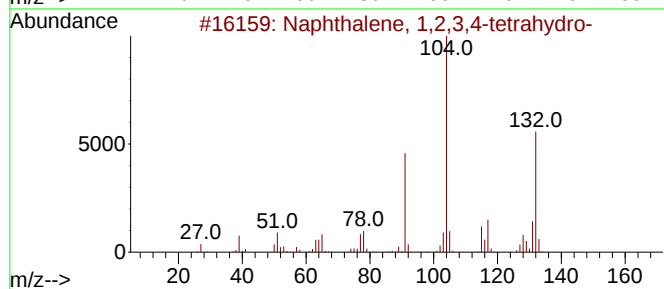
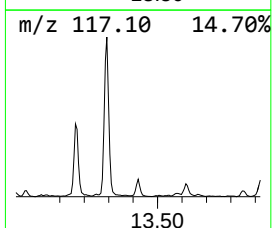
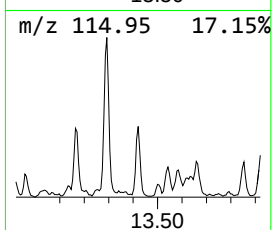
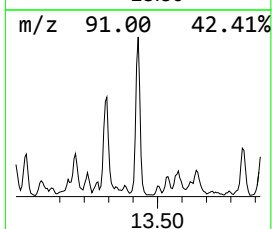
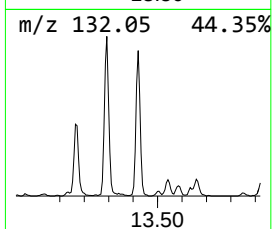
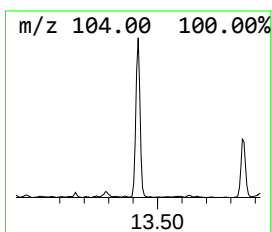
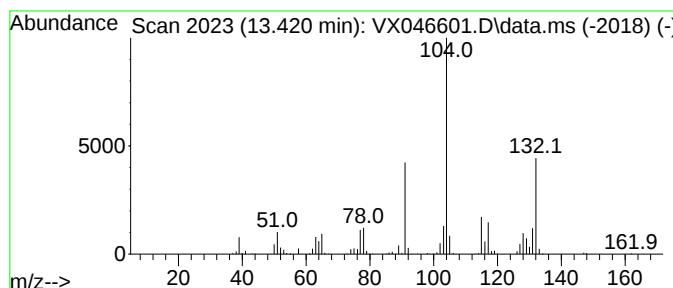
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	14.32 ug/l	217564	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	53
4			N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	50
5			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

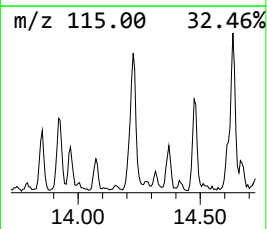
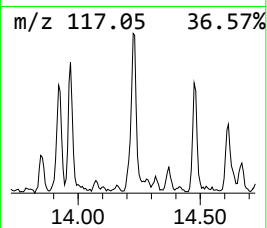
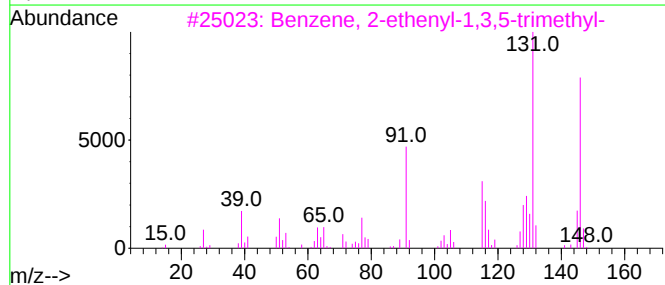
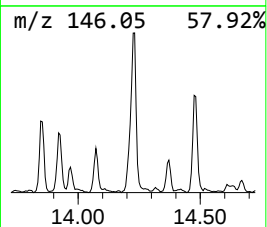
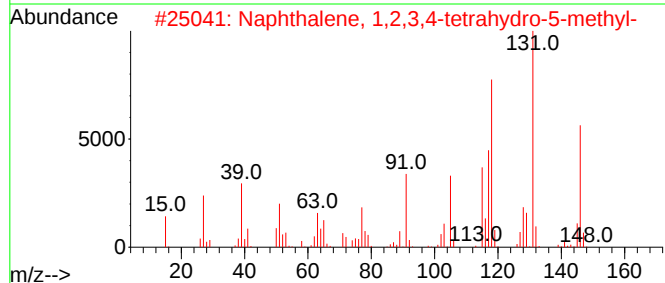
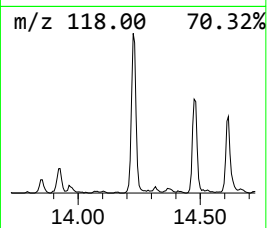
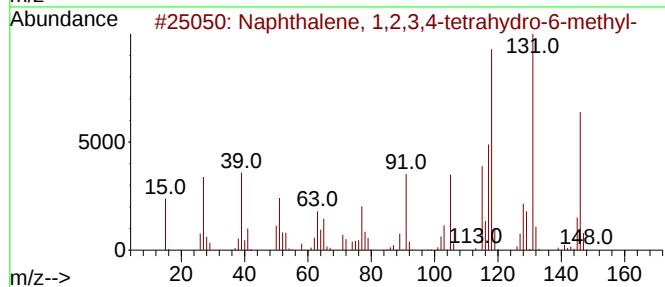
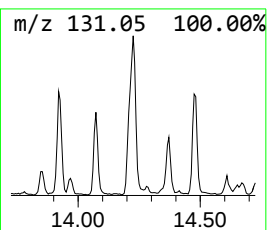
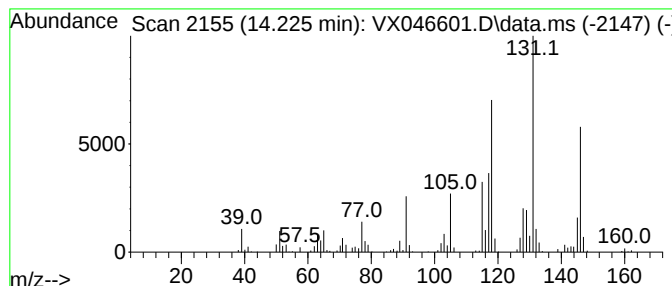
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	16.97 ug/l	257845	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
5			1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GW-MW901-060425

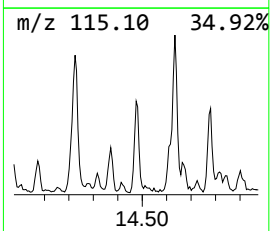
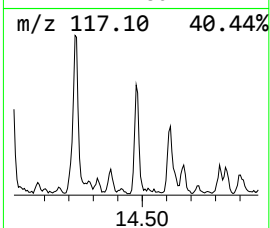
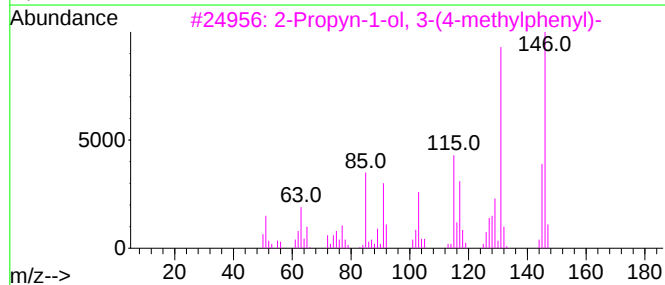
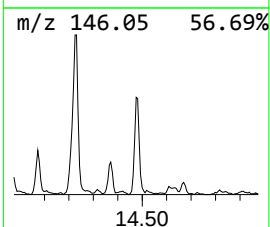
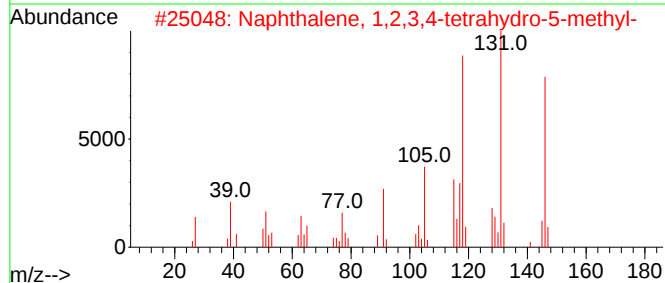
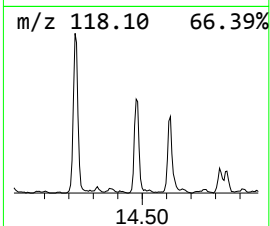
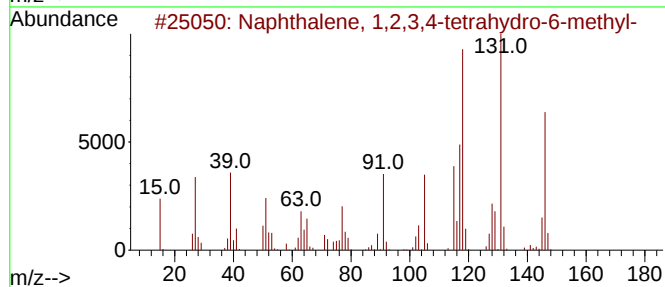
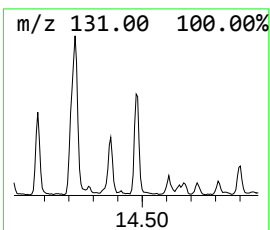
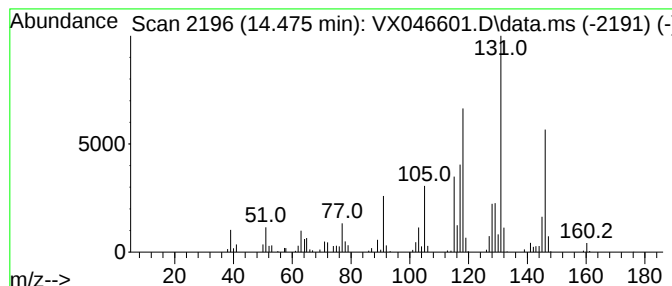
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.475	9.09 ug/l	138097	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	94
3			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	76
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061025\
Data File : VX046601.D
Acq On : 10 Jun 2025 14:38
Operator : JC/MD
Sample : Q2230-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GW-MW901-060425

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.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
Benzene, 1-ethy...	11.378	22.8	ug/l	345847	4	12.018	759761	50.0
Benzene, 1-ethy...	11.610	15.7	ug/l	239196	4	12.018	759761	50.0
Benzene, 1,2,3-...	12.085	19.2	ug/l	292222	4	12.018	759761	50.0
Indane	12.244	13.2	ug/l	200586	4	12.018	759761	50.0
o-Cymene	12.317	10.4	ug/l	157984	4	12.018	759761	50.0
Indan, 1-methyl-	12.695	12.9	ug/l	195774	4	12.018	759761	50.0
Benzene, 1-ethe...	13.292	19.1	ug/l	289968	4	12.018	759761	50.0
Naphthalene, 1,...	13.420	14.3	ug/l	217564	4	12.018	759761	50.0
Naphthalene, 1,...	14.225	17.0	ug/l	257845	4	12.018	759761	50.0
Naphthalene, 1,...	14.475	9.1	ug/l	138097	4	12.018	759761	50.0