

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
 Data File : VX044045.D
 Acq On : 27 Nov 2024 16:57
 Operator : JC/MD
 Sample : P5021-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-06

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\82X112124W.M
 Title : SW846 8260

Signal : TIC: VX044045.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.587	66	82	83	rBV4	16405	53278	3.07%	0.274%
2	2.770	269	276	279	rBV3	13752	28985	1.67%	0.149%
3	2.812	279	283	288	rVB2	14146	26787	1.54%	0.138%
4	3.087	322	328	340	rVB3	24011	57846	3.33%	0.297%
5	3.721	424	432	441	rBV2	15544	38059	2.19%	0.196%
6	4.202	500	511	518	rBV2	19418	57533	3.31%	0.296%
7	4.294	518	526	537	rVB3	39841	113433	6.53%	0.583%
8	5.379	694	704	713	rBV	97121	291416	16.78%	1.498%
9	5.544	714	731	737	rVV2	153289	506800	29.18%	2.604%
10	5.605	738	741	758	rVB2	74331	217628	12.53%	1.118%
11	5.818	765	776	788	rVB3	21780	66658	3.84%	0.343%
12	5.952	788	798	805	rBV	117632	311985	17.96%	1.603%
13	6.031	805	811	823	rVV	51054	142673	8.21%	0.733%
14	6.159	823	832	834	rVV4	14628	42916	2.47%	0.221%
15	6.239	835	845	857	rVV	139105	448240	25.80%	2.303%
16	6.342	857	862	872	rVB5	12645	36832	2.12%	0.189%
17	6.757	921	930	940	rBV	253575	638908	36.78%	3.283%
18	7.165	988	997	1005	rVB3	17073	43070	2.48%	0.221%
19	7.366	1020	1030	1044	rVB6	38852	120503	6.94%	0.619%
20	7.495	1044	1051	1055	rBV3	16076	34790	2.00%	0.179%
21	7.555	1055	1061	1064	rVV3	16002	35042	2.02%	0.180%
22	7.598	1065	1068	1073	rVV	13261	25233	1.45%	0.130%
23	7.653	1073	1077	1084	rVB5	11697	21448	1.23%	0.110%
24	7.903	1111	1118	1123	rVV5	21539	55546	3.20%	0.285%
25	7.976	1123	1130	1140	rVB	129973	265663	15.29%	1.365%
26	8.104	1142	1151	1161	rBV2	194963	448980	25.85%	2.307%
27	8.183	1161	1164	1170	rVB2	16445	25841	1.49%	0.133%
28	8.427	1199	1204	1211	rBV	74697	134320	7.73%	0.690%
29	8.500	1211	1216	1224	rVB3	30556	58486	3.37%	0.301%
30	8.580	1224	1229	1231	rBV2	21410	35550	2.05%	0.183%
31	8.647	1235	1240	1247	rVV	516885	817757	47.08%	4.202%
32	8.720	1248	1252	1258	rVB2	19554	32958	1.90%	0.169%
33	8.811	1263	1267	1278	rVB2	16207	32099	1.85%	0.165%
34	9.098	1306	1314	1319	rVB3	11711	22751	1.31%	0.117%
35	9.159	1319	1324	1333	rVV2	27904	52443	3.02%	0.269%
36	9.244	1333	1338	1344	rVB	25449	39587	2.28%	0.203%

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37	9.451	1364	1372	1377	rVV4	31571	66689	3.84%	0.343%
38	9.945	1448	1453	1460	rBV	76334	111121	6.40%	0.571%
39	10.012	1460	1464	1465	rVV2	17055	23332	1.34%	0.120%
40	10.049	1465	1470	1475	rVV	508901	717493	41.30%	3.687%
41	10.092	1475	1477	1484	rVB4	21900	37634	2.17%	0.193%
42	10.214	1490	1497	1500	rBV2	78634	137258	7.90%	0.705%
43	10.244	1500	1502	1507	rVV	36491	41464	2.39%	0.213%
44	10.299	1507	1511	1517	rVB	50001	76537	4.41%	0.393%
45	10.384	1517	1525	1529	rBV2	63714	112207	6.46%	0.577%
46	10.640	1563	1567	1574	rVB2	27188	37216	2.14%	0.191%
47	10.793	1582	1592	1599	rBV10	10605	34703	2.00%	0.178%
48	10.896	1599	1609	1614	rBV2	38678	77970	4.49%	0.401%
49	10.957	1614	1619	1625	rBV	311500	426249	24.54%	2.190%
50	11.012	1625	1628	1634	rVB	28289	37203	2.14%	0.191%
51	11.079	1634	1639	1649	rBV	423190	598084	34.43%	3.073%
52	11.158	1649	1652	1660	rVB4	18373	40087	2.31%	0.206%
53	11.305	1671	1676	1683	rBV	665827	864037	49.74%	4.440%
54	11.451	1695	1700	1703	rVV4	17177	26804	1.54%	0.138%
55	11.494	1703	1707	1712	rVB2	22476	34540	1.99%	0.177%
56	11.610	1721	1726	1730	rBV	67672	88287	5.08%	0.454%
57	11.750	1745	1749	1755	rVB2	34076	51835	2.98%	0.266%
58	11.866	1761	1768	1769	rBV	41456	51160	2.95%	0.263%
59	11.890	1769	1772	1776	rVB	50561	60546	3.49%	0.311%
60	12.018	1788	1793	1800	rVV	513051	645153	37.14%	3.315%
61	12.085	1800	1804	1808	rVB	40130	47748	2.75%	0.245%
62	12.177	1816	1819	1822	rVB	24043	26673	1.54%	0.137%
63	12.244	1824	1830	1837	rVV	1374850	1737101	100.00%	8.927%
64	12.311	1837	1841	1851	rVV2	164357	249690	14.37%	1.283%
65	12.402	1851	1856	1861	rVV2	132266	165291	9.52%	0.849%
66	12.469	1861	1867	1872	rVB3	53535	93349	5.37%	0.480%
67	12.530	1872	1877	1884	rBV5	17526	31361	1.81%	0.161%
68	12.609	1885	1890	1894	rBV	1123398	1334196	76.81%	6.856%
69	12.640	1894	1895	1899	rVV	54947	56009	3.22%	0.288%
70	12.695	1899	1904	1912	rVB2	335688	539824	31.08%	2.774%
71	12.835	1923	1927	1932	rVB3	22490	33337	1.92%	0.171%
72	12.920	1932	1941	1947	rBV	692390	863884	49.73%	4.439%
73	13.024	1953	1958	1966	rVB3	57262	85494	4.92%	0.439%
74	13.115	1966	1973	1975	rVV3	40806	73642	4.24%	0.378%
75	13.164	1978	1981	1989	rVV	210879	283179	16.30%	1.455%
76	13.237	1989	1993	1994	rVV2	26111	36177	2.08%	0.186%

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77	13.256	1994	1996	1998	rVV	51008	64863	3.73%	0.333%
78	13.292	1998	2002	2007	rVV2	445952	665806	38.33%	3.422%
79	13.341	2007	2010	2012	rVV2	59197	75684	4.36%	0.389%
80	13.365	2012	2014	2020	rVV	55228	81216	4.68%	0.417%
81	13.420	2020	2023	2028	rVB2	51287	66420	3.82%	0.341%
82	13.506	2032	2037	2039	rVV2	22234	32491	1.87%	0.167%
83	13.542	2039	2043	2046	rVV	96011	139812	8.05%	0.718%
84	13.573	2046	2048	2053	rVV3	134631	185061	10.65%	0.951%
85	13.640	2053	2059	2060	rVV2	49612	79080	4.55%	0.406%
86	13.670	2060	2064	2069	rVB2	83870	145512	8.38%	0.748%
87	13.774	2074	2081	2091	rVV	1047312	1337737	77.01%	6.875%
88	13.926	2099	2106	2110	rVV2	35426	57262	3.30%	0.294%
89	13.969	2110	2113	2122	rVB	31901	47753	2.75%	0.245%
90	14.073	2125	2130	2136	rBV	71725	92852	5.35%	0.477%
91	14.213	2149	2153	2159	rBV	58567	91574	5.27%	0.471%
92	14.316	2166	2170	2175	rBV	100896	126174	7.26%	0.648%
93	14.371	2175	2179	2183	rVV	43910	57403	3.30%	0.295%
94	14.597	2211	2216	2219	rBV3	17722	27875	1.60%	0.143%
95	14.633	2219	2222	2227	rVB	104485	131431	7.57%	0.675%
96	14.780	2241	2246	2255	rVB	218384	296203	17.05%	1.522%
97	15.615	2376	2383	2386	rBV	24923	41183	2.37%	0.212%
98	15.834	2417	2419	2429	rVB5	11257	20479	1.18%	0.105%
99	16.322	2492	2499	2507	rBV	71402	130439	7.51%	0.670%
100	16.688	2551	2559	2570	rVB2	12975	28471	1.64%	0.146%

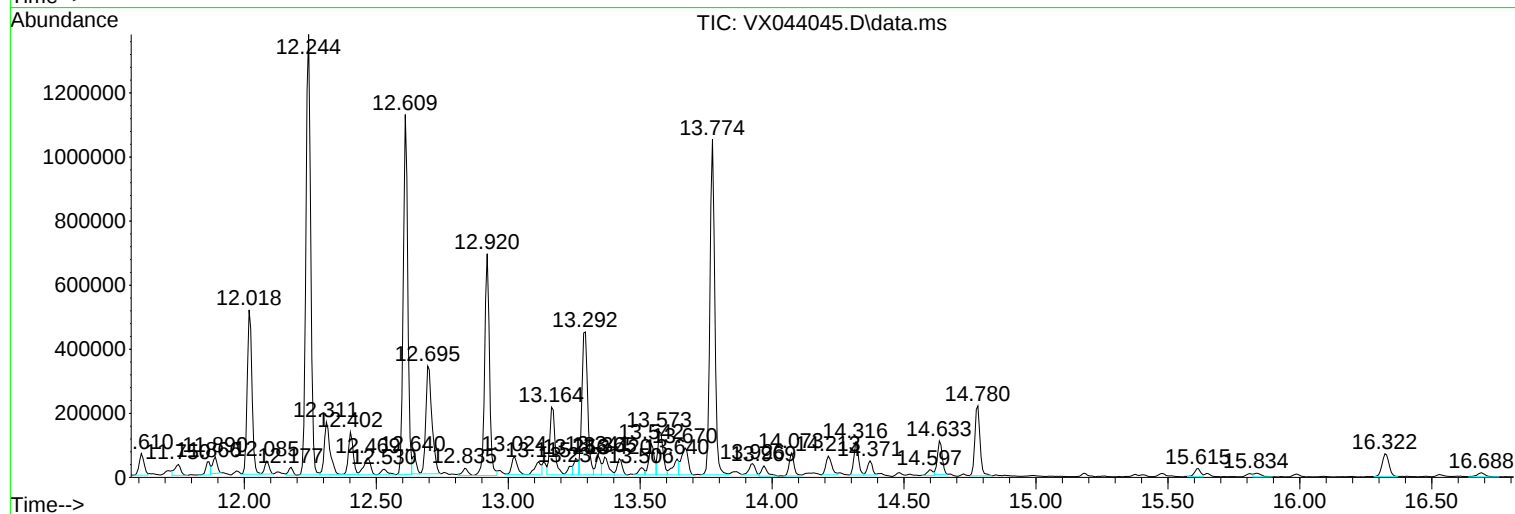
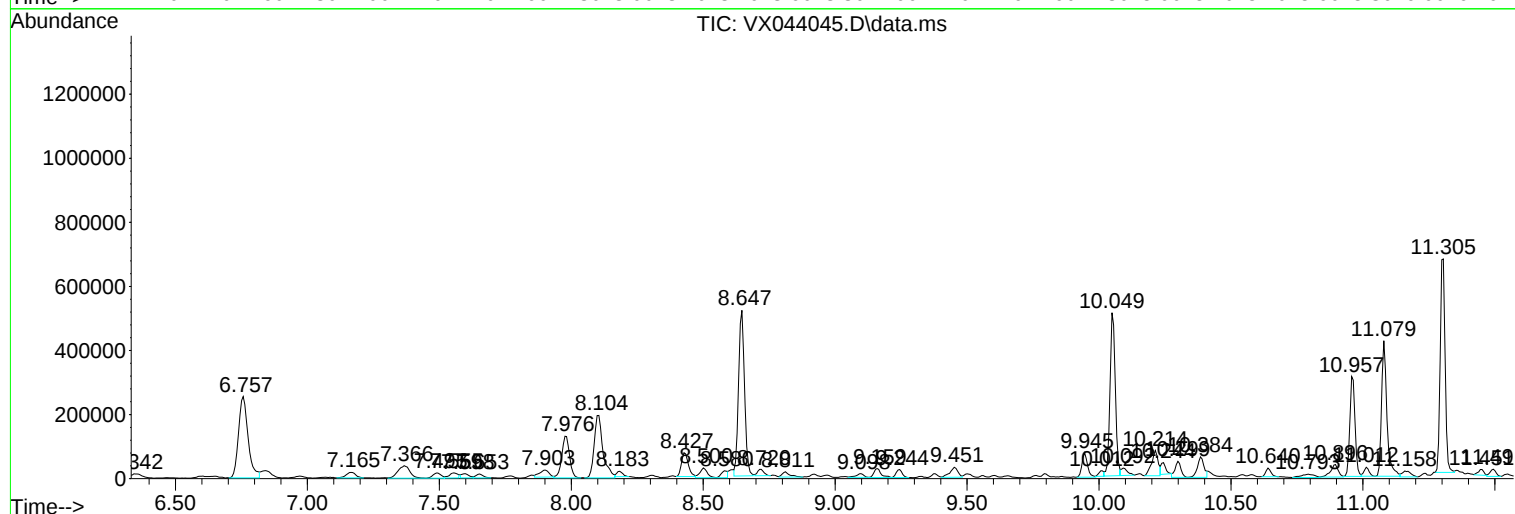
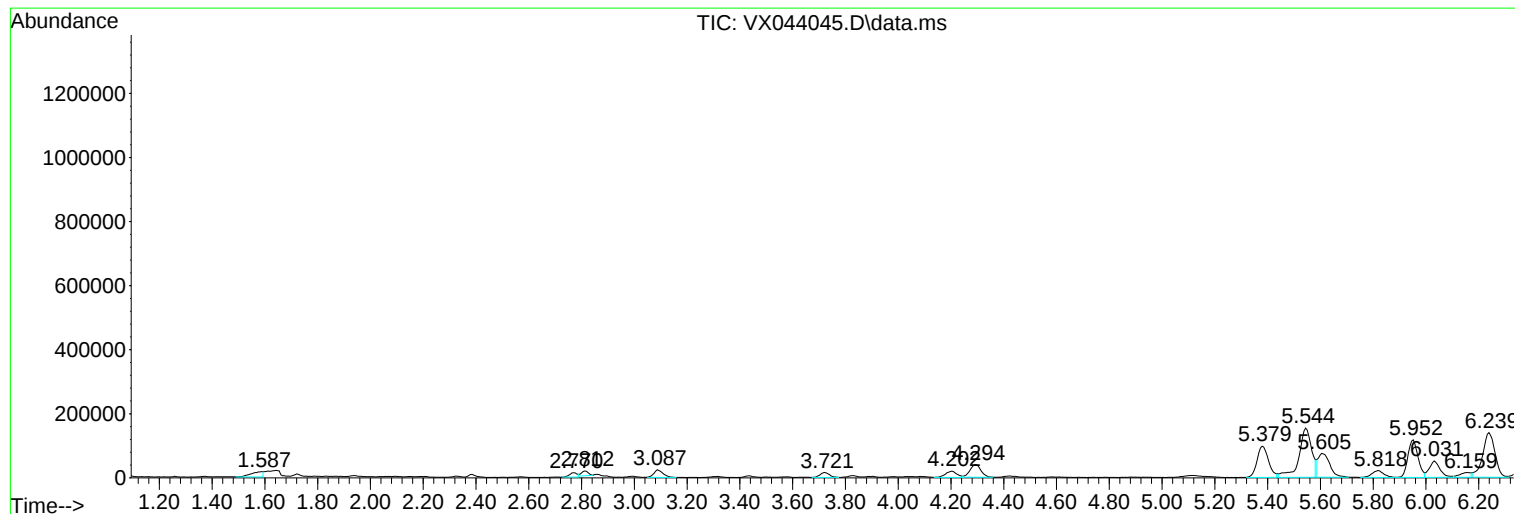
Sum of corrected areas: 19459389

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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



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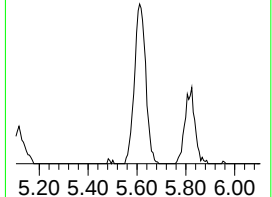
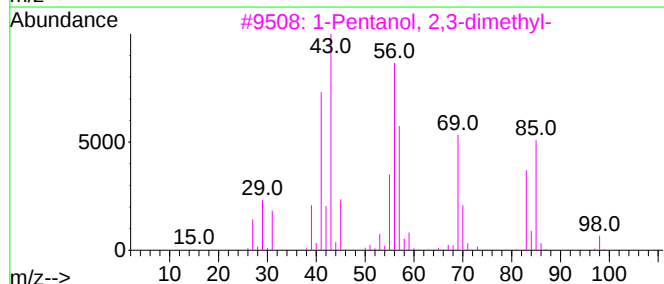
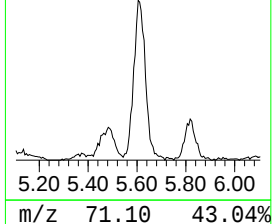
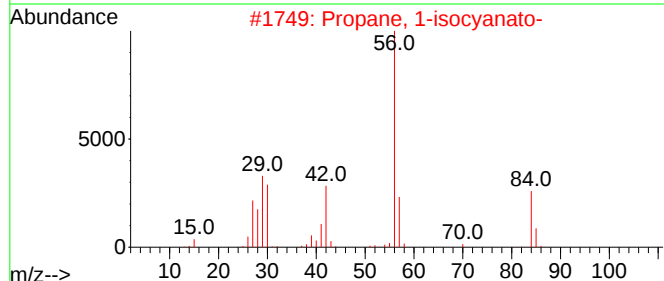
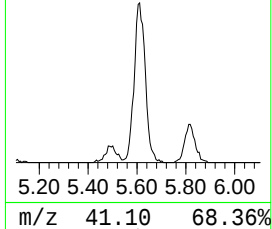
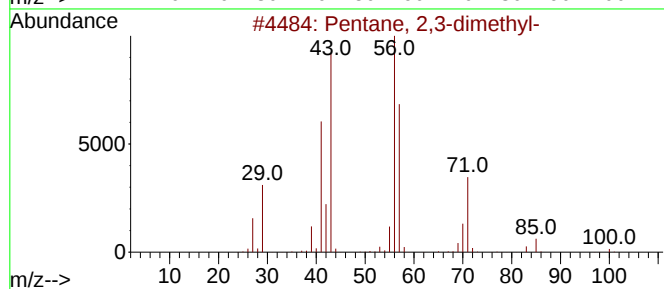
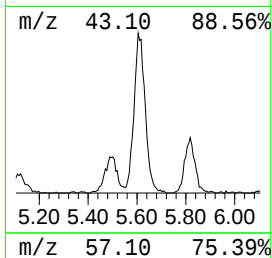
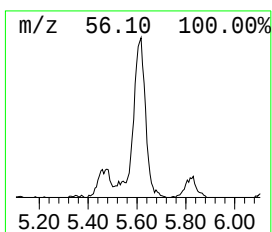
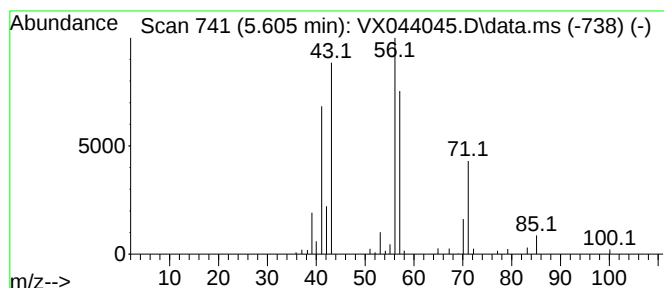
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.605	21.47 ug/l	217628	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	43
3			1-Pentanol, 2,3-dimethyl-	116	C7H16O	010143-23-4	42
4			Ethane, isocyanato-	71	C3H5NO	000109-90-0	40
5			1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	38



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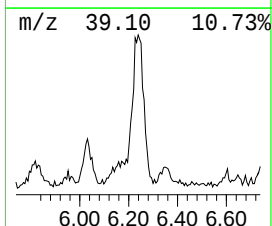
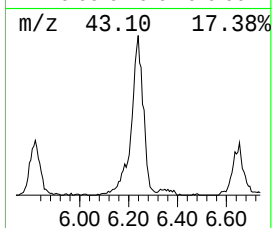
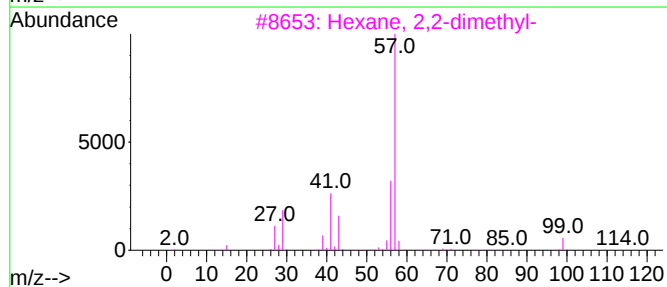
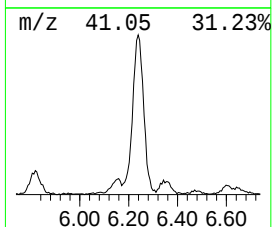
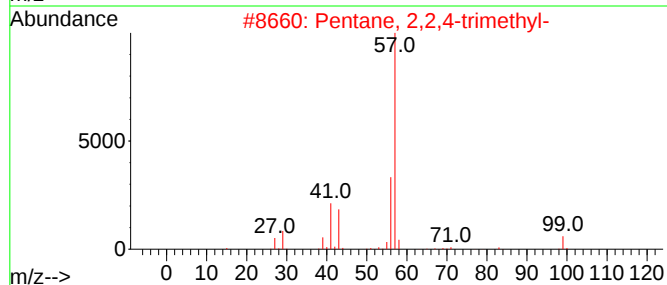
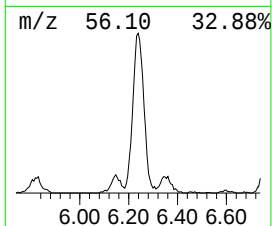
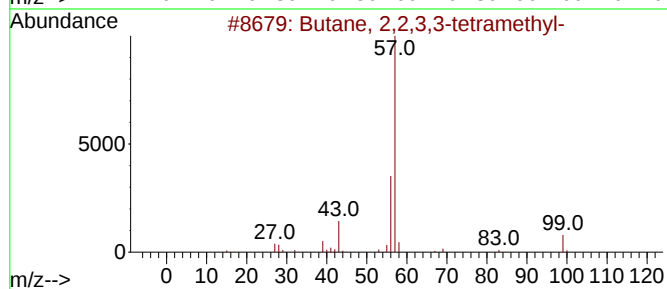
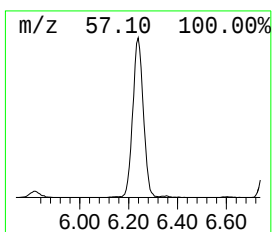
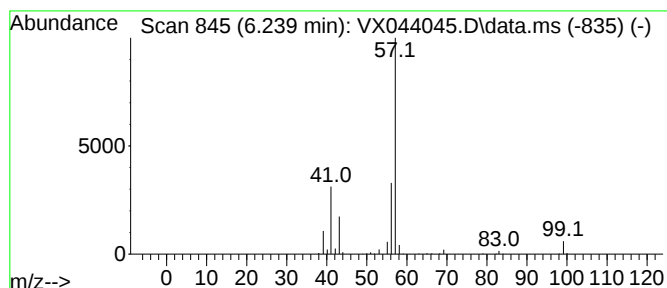
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2,2,3,3-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.239	35.08 ug/l	448240	1,4-Difluorobenzene	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64



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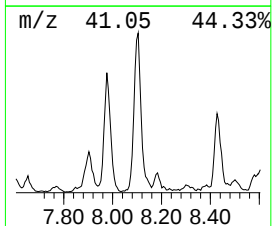
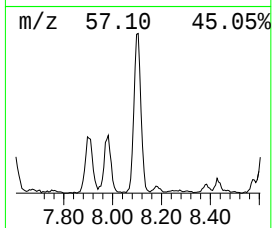
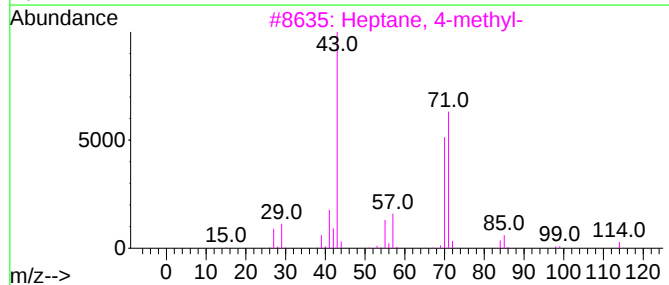
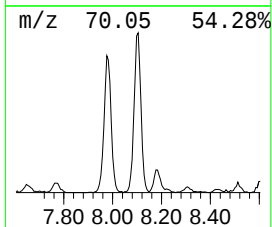
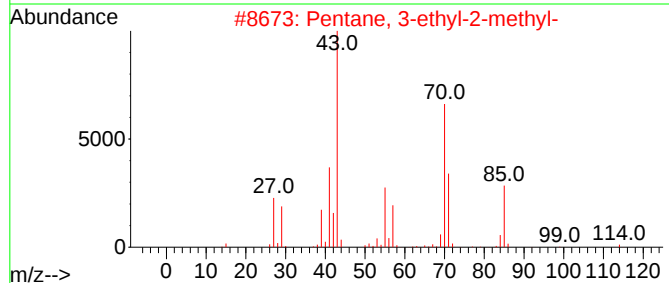
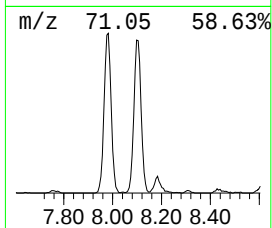
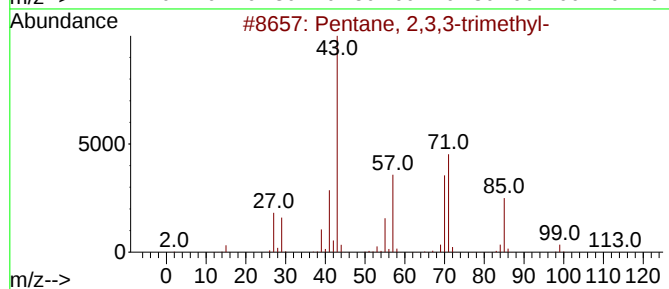
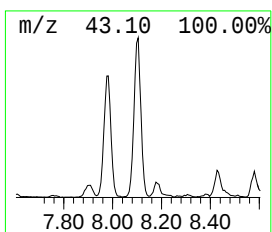
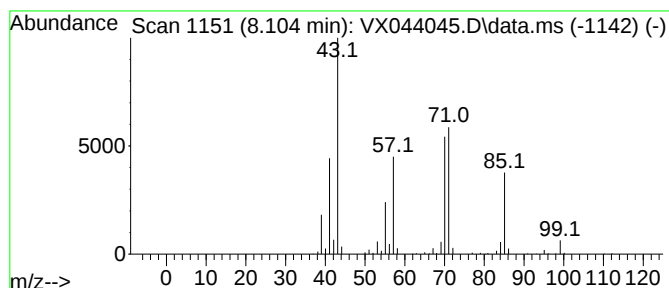
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	35.14 ug/l	448980	1,4-Difluorobenzene	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2			Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	64
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	59
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

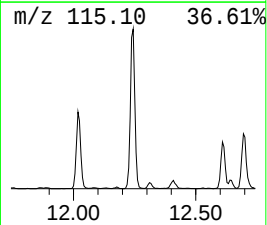
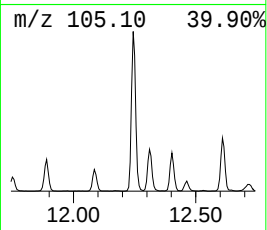
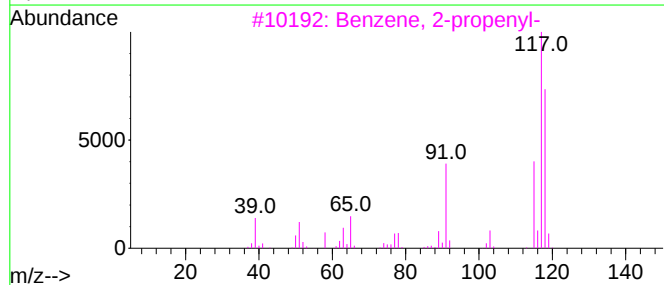
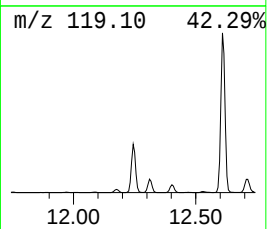
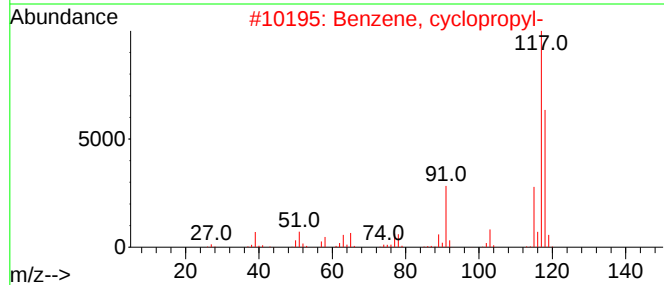
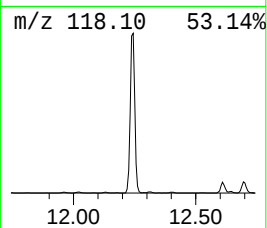
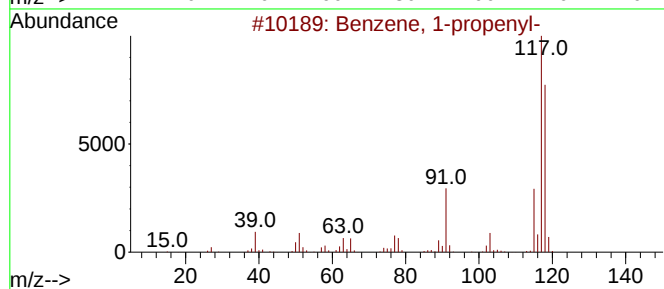
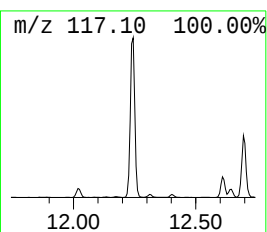
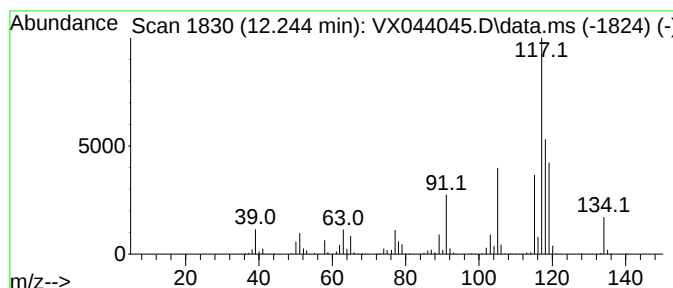
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	134.63 ug/l	1737100	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4			Indane	118	C9H10	000496-11-7	60
5			Deltacyclene	118	C9H10	007785-10-6	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

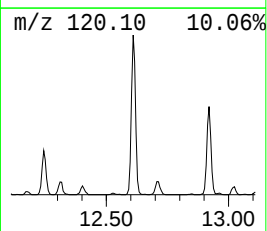
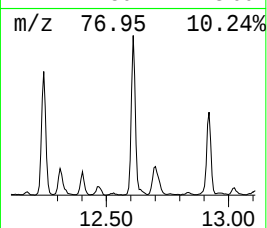
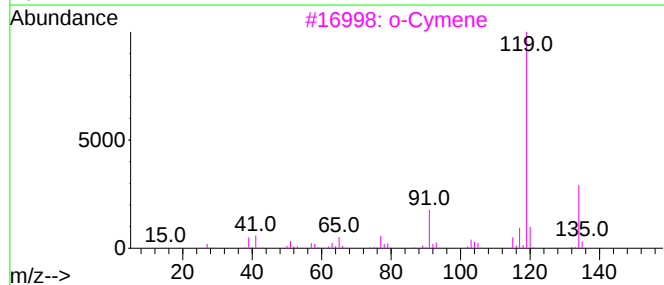
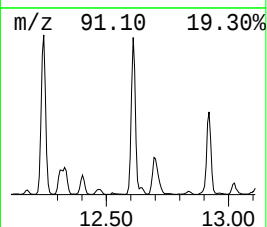
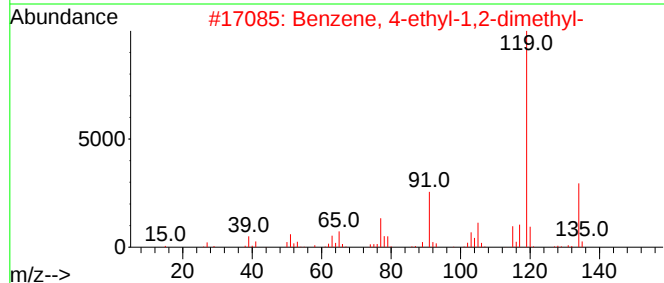
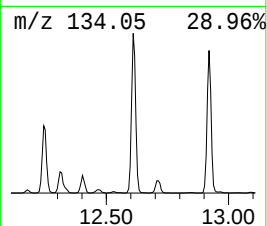
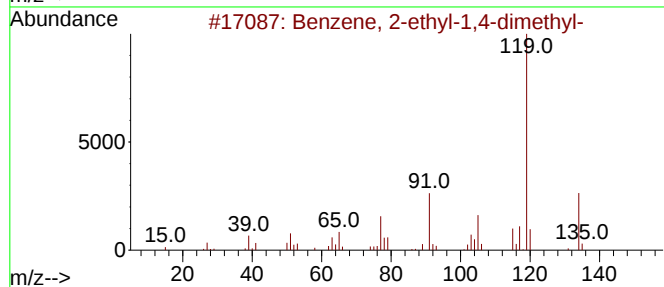
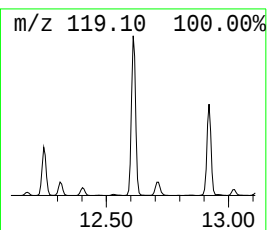
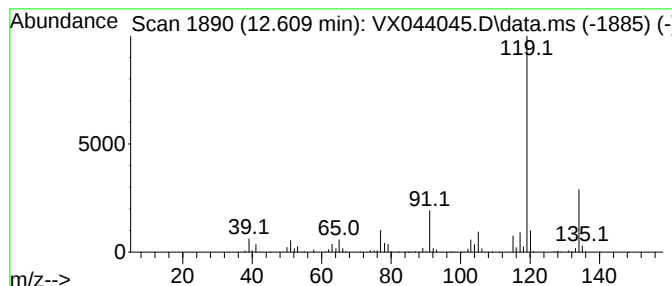
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.609	103.40 ug/l	1334200	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

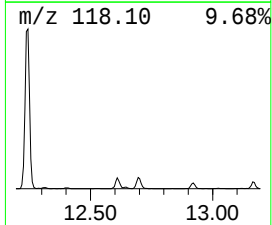
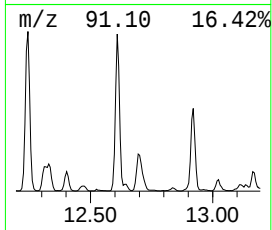
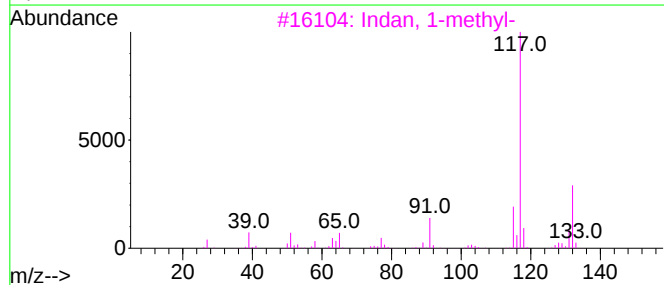
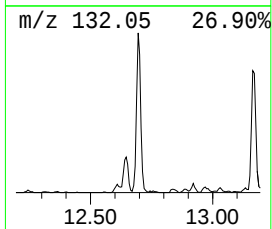
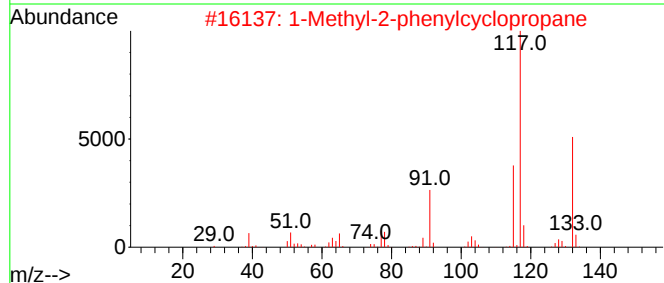
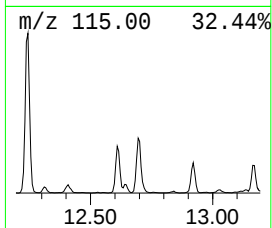
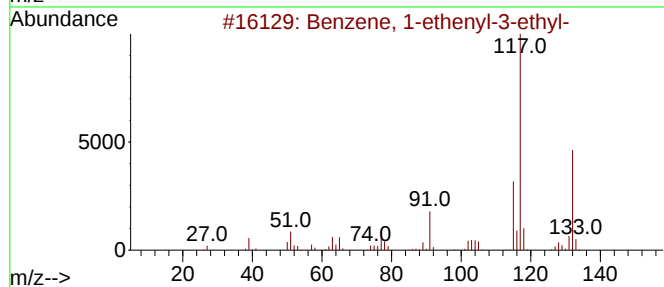
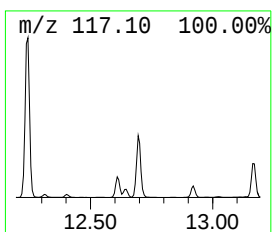
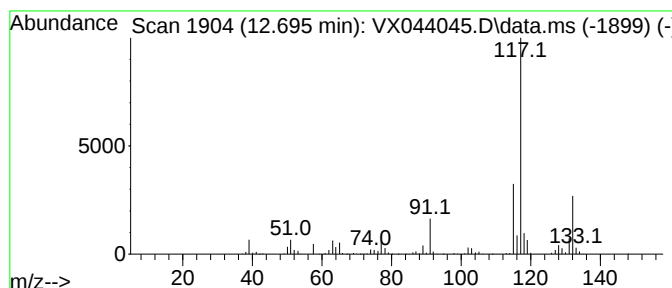
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	41.84 ug/l	539824	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	90
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
3			Indan, 1-methyl-	132	C10H12	000767-58-8	87
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

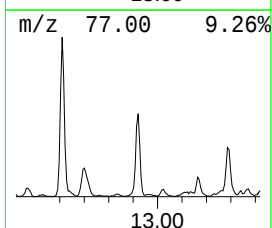
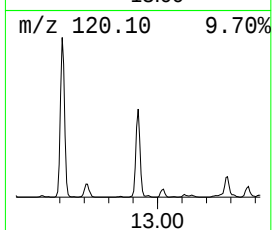
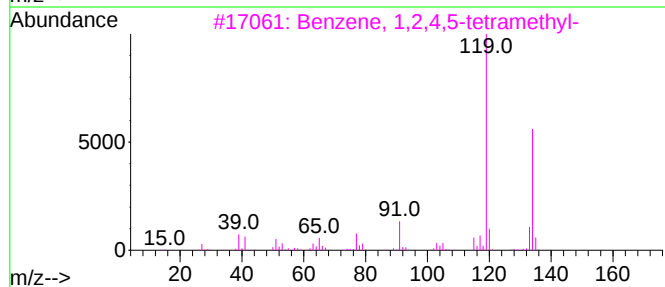
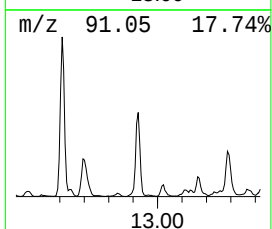
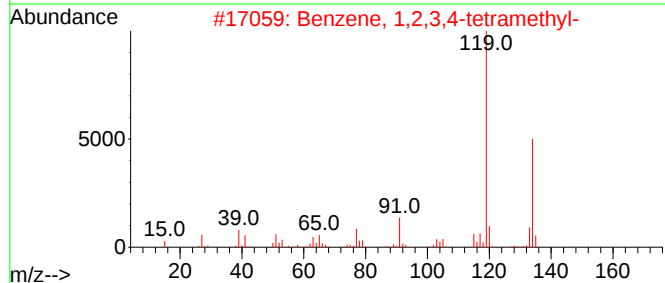
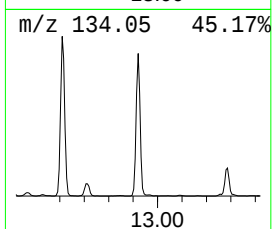
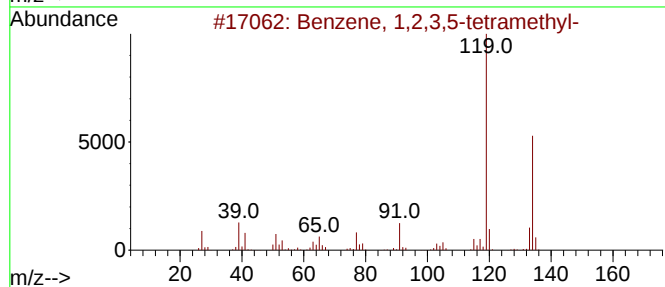
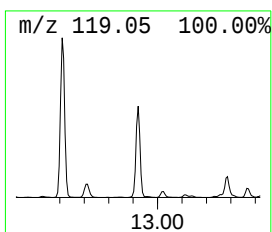
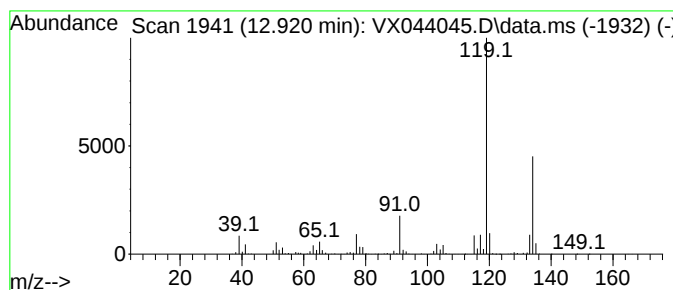
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	66.95 ug/l	863884	1,4-Dichlorobenzene-d4	12.018

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

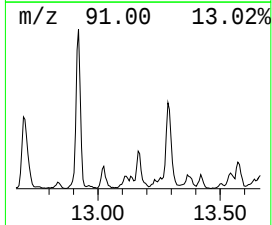
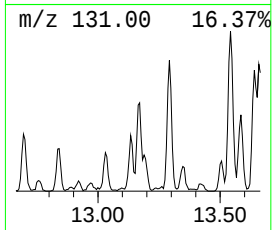
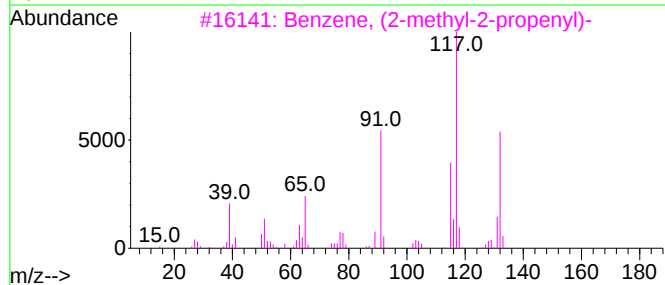
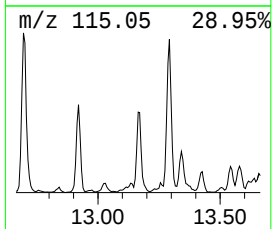
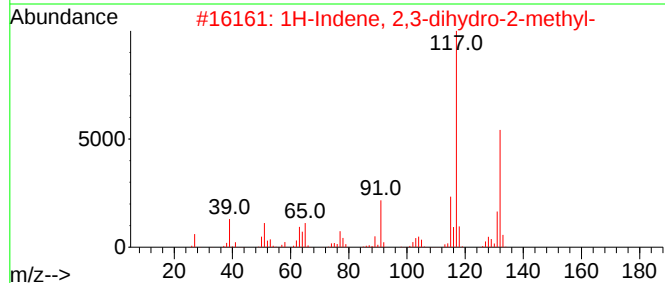
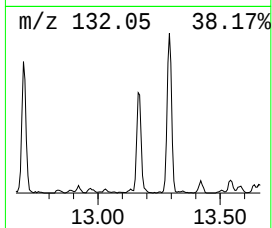
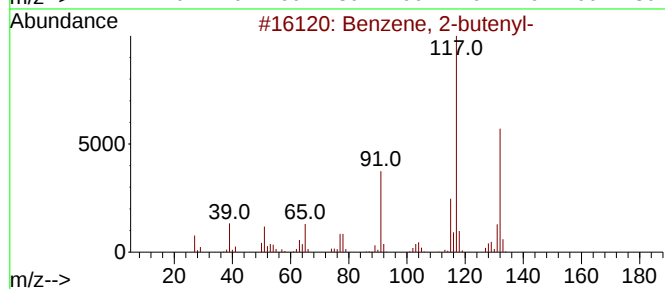
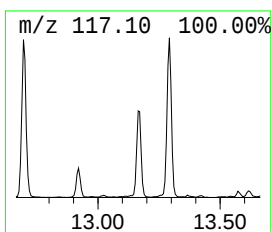
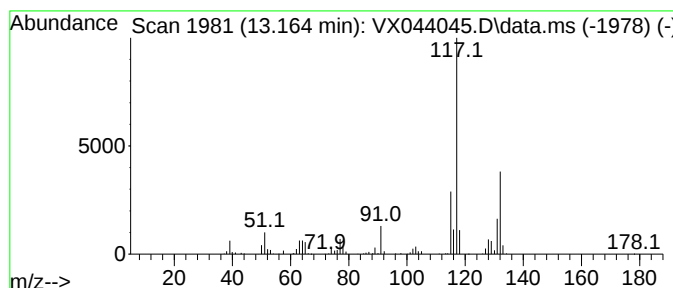
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-butenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.164	21.95 ug/l	283179	1,4-Dichlorobenzene-d4	12.018

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
2			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
3			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
5			Indan, 1-methyl-	132	C10H12	000767-58-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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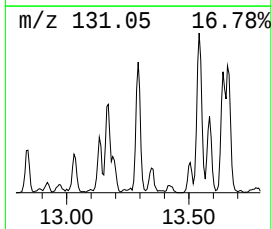
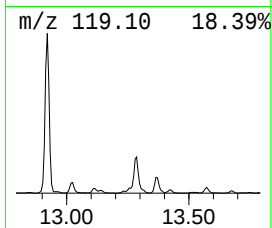
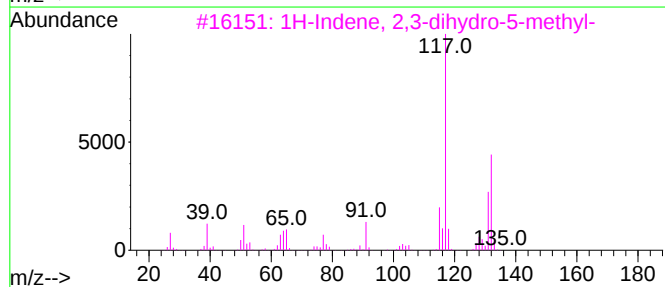
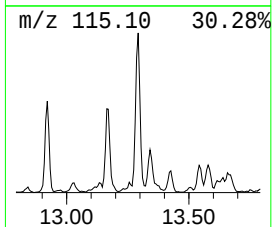
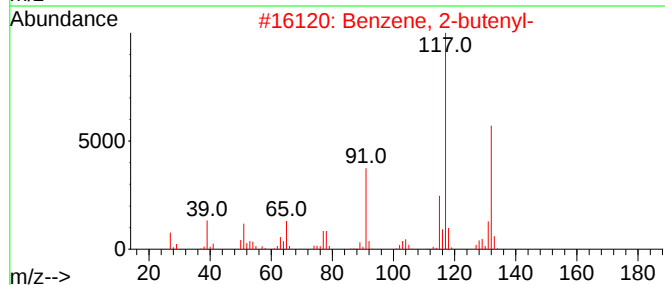
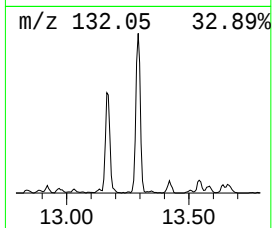
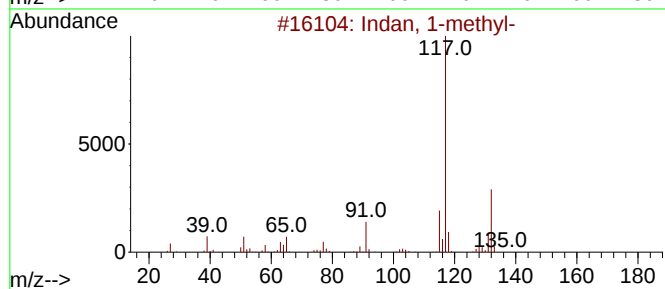
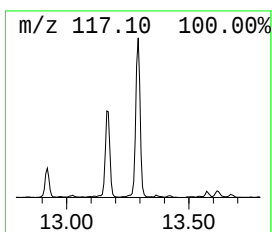
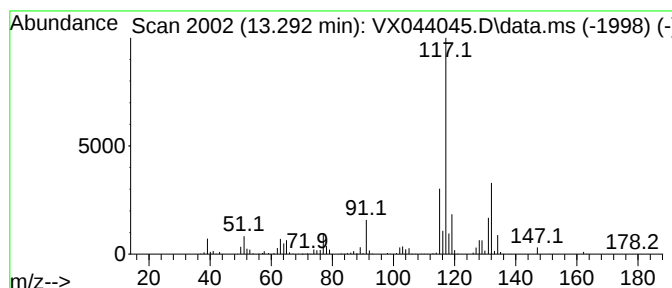
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Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	51.60 ug/l	665806	1,4-Dichlorobenzene-d4	12.018

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	83
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

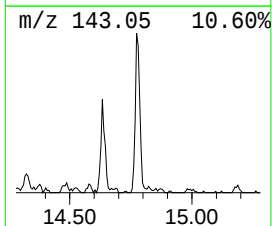
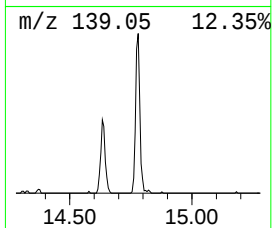
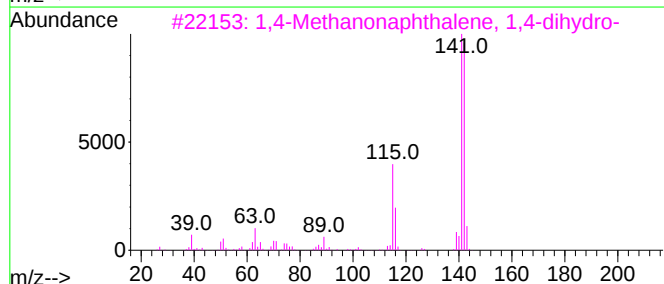
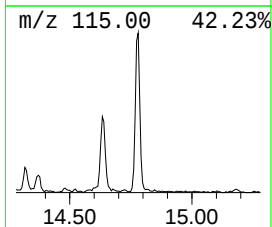
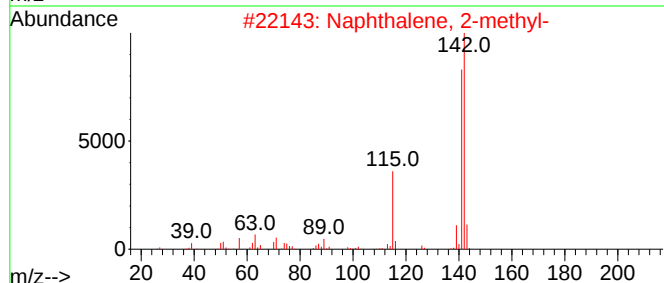
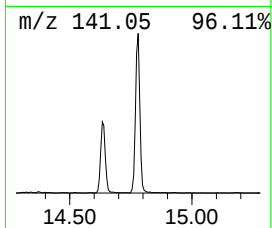
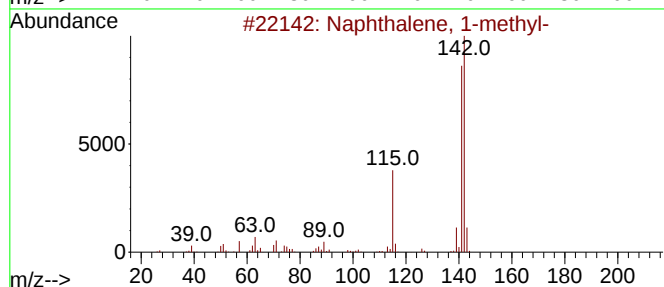
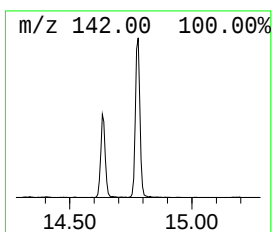
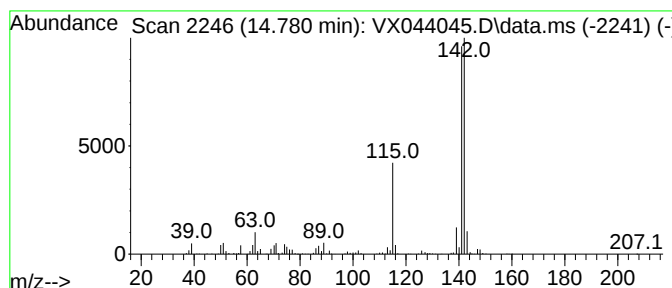
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	22.96 ug/l	296203	1,4-Dichlorobenzene-d4	12.018

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
Data File : VX044045.D
Acq On : 27 Nov 2024 16:57
Operator : JC/MD
Sample : P5021-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,3-di...	5.605	21.5	ug/l	217628	1	5.544	506800	50.0
Butane, 2,2,3,3...	6.239	35.1	ug/l	448240	2	6.757	638908	50.0
Pentane, 2,3,3-...	8.104	35.1	ug/l	448980	2	6.757	638908	50.0
Benzene, 1-prop...	12.244	134.6	ug/l	1737100	4	12.018	645153	50.0
Benzene, 2-ethy...	12.609	103.4	ug/l	1334200	4	12.018	645153	50.0
Benzene, 1-ethe...	12.695	41.8	ug/l	539824	4	12.018	645153	50.0
Benzene, 1,2,3,...	12.920	67.0	ug/l	863884	4	12.018	645153	50.0
Benzene, 2-bute...	13.164	21.9	ug/l	283179	4	12.018	645153	50.0
Indan, 1-methyl-	13.292	51.6	ug/l	665806	4	12.018	645153	50.0
Naphthalene, 1-...	14.780	23.0	ug/l	296203	4	12.018	645153	50.0