

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
 Data File : VX033304.D
 Acq On : 08 Dec 2022 00:10
 Operator : JC/MD
 Sample : N5906-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-13R

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\82X120722W.M
 Title : SW846 8260

Signal : TIC: VX033304.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.368	41	46	52	rBV	157740	171192	2.35%	0.190%
2	1.752	103	109	121	rVB	608266	817878	11.25%	0.907%
3	1.959	138	143	156	rVB3	237557	448384	6.17%	0.498%
4	2.069	156	161	168	rVV	143902	199129	2.74%	0.221%
5	2.166	172	177	181	rVV	76297	114428	1.57%	0.127%
6	2.215	181	185	198	rVV	276103	415171	5.71%	0.461%
7	2.782	259	278	281	rBV2	183488	496838	6.83%	0.551%
8	2.825	281	285	289	rVV	261636	561430	7.72%	0.623%
9	2.867	289	292	304	rVV2	240016	584072	8.03%	0.648%
10	2.983	304	311	322	rVV2	31669	106325	1.46%	0.118%
11	3.105	322	331	349	rVB3	334176	792714	10.90%	0.880%
12	3.319	356	366	378	rBV	58013	130845	1.80%	0.145%
13	3.447	378	387	400	rVB	90812	198794	2.73%	0.221%
14	3.733	427	434	443	rVV	181788	457010	6.28%	0.507%
15	3.837	443	451	458	rVV2	119866	318666	4.38%	0.354%
16	3.910	458	463	472	rVV2	58596	190988	2.63%	0.212%
17	4.105	486	495	504	rVV	74031	205809	2.83%	0.228%
18	4.215	504	513	518	rVV2	38225	113155	1.56%	0.126%
19	4.306	518	528	540	rVV	589108	1677194	23.06%	1.861%
20	4.428	540	548	565	rVB	59682	179668	2.47%	0.199%
21	5.196	663	674	690	rVB2	157687	507717	6.98%	0.563%
22	5.385	690	705	711	rBV3	71749	224403	3.09%	0.249%
23	5.477	711	720	727	rVV2	227834	808032	11.11%	0.897%
24	5.556	728	733	739	rVV	164093	497094	6.84%	0.552%
25	5.623	740	744	765	rVB4	113370	370709	5.10%	0.411%
26	5.830	765	778	791	rBV3	97358	298628	4.11%	0.331%
27	5.958	791	799	803	rVV2	45808	110817	1.52%	0.123%
28	6.044	803	813	823	rVV	1129481	2998460	41.23%	3.327%
29	6.257	824	848	857	rVV4	259807	1311376	18.03%	1.455%
30	6.361	857	865	876	rVB2	84355	251639	3.46%	0.279%
31	6.611	893	906	914	rBV2	54237	141274	1.94%	0.157%
32	6.763	923	931	937	rBV	234564	612689	8.42%	0.680%
33	6.848	941	945	959	rVV4	161592	421463	5.80%	0.468%
34	7.171	989	998	1007	rBV	133569	302071	4.15%	0.335%
35	7.379	1019	1032	1045	rBV5	355417	966841	13.29%	1.073%
36	7.659	1073	1078	1090	rVB	83874	159786	2.20%	0.177%

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37	7.988	1122	1132	1143	rVB	182915	408486	5.62%	0.453%
38	8.110	1143	1152	1155	rBV	286720	619033	8.51%	0.687%
39	8.141	1155	1157	1162	rVV	320405	501982	6.90%	0.557%
40	8.190	1162	1165	1175	rVV3	88502	172548	2.37%	0.191%

41	8.507	1211	1217	1226	rVB2	201800	369361	5.08%	0.410%
42	8.610	1226	1234	1236	rBV4	68000	137419	1.89%	0.152%
43	8.647	1236	1240	1246	rVB	204735	341876	4.70%	0.379%
44	8.720	1246	1252	1258	rBV	349593	534564	7.35%	0.593%
45	8.958	1287	1291	1300	rVV2	169814	311425	4.28%	0.346%

46	9.049	1300	1306	1311	rVV	146282	240181	3.30%	0.266%
47	9.159	1320	1324	1330	rBV	114813	173253	2.38%	0.192%
48	10.055	1466	1471	1476	rVB	480478	610020	8.39%	0.677%
49	10.195	1489	1494	1500	rVB	5350115	6802018	93.53%	7.547%
50	10.305	1506	1512	1518	rVB	2588719	3219407	44.27%	3.572%

51	10.640	1562	1567	1578	rVB	835345	1059375	14.57%	1.175%
52	10.963	1614	1620	1629	rBV	1429408	1760881	24.21%	1.954%
53	11.079	1635	1639	1644	rBV	344705	440183	6.05%	0.488%
54	11.305	1667	1676	1683	rBV	4023795	4754013	65.37%	5.275%
55	11.402	1683	1692	1696	rVV	1504624	2107936	28.98%	2.339%

56	11.451	1696	1700	1708	rVB	249451	305911	4.21%	0.339%
57	11.616	1721	1727	1737	rBV	2856684	3644983	50.12%	4.044%
58	11.750	1744	1749	1755	rVB	2936849	3590832	49.37%	3.984%
59	11.866	1763	1768	1770	rBV	179330	231379	3.18%	0.257%
60	11.890	1770	1772	1778	rVB	244395	266212	3.66%	0.295%

61	11.969	1778	1785	1789	rBV	111709	137137	1.89%	0.152%
62	12.024	1789	1794	1798	rVV	544035	718482	9.88%	0.797%
63	12.085	1799	1804	1809	rVB	1784009	2198099	30.22%	2.439%
64	12.244	1824	1830	1837	rBV	6007349	7272721	100.00%	8.070%
65	12.311	1837	1841	1852	rVB2	678071	1197446	16.46%	1.329%

66	12.408	1852	1857	1862	rBV2	284660	375035	5.16%	0.416%
67	12.463	1862	1866	1872	rVB	562030	649221	8.93%	0.720%
68	12.530	1872	1877	1880	rBV	451972	608547	8.37%	0.675%
69	12.555	1880	1881	1885	rVB	238824	193802	2.66%	0.215%
70	12.616	1885	1891	1894	rBV	2671202	3235306	44.49%	3.590%

71	12.646	1894	1896	1900	rVB	471759	483263	6.64%	0.536%
72	12.701	1900	1905	1912	rBV	1515726	2037372	28.01%	2.261%
73	12.847	1924	1929	1934	rVB2	284448	381060	5.24%	0.423%
74	12.920	1934	1941	1945	rBV	1615148	1978899	27.21%	2.196%
75	12.963	1945	1948	1954	rVB	1076796	1236968	17.01%	1.372%

76	13.024	1954	1958	1965	rBV3	126606	203962	2.80%	0.226%
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77	13.170	1978	1982	1987	rVB	1277268	1499881	20.62%	1.664%
78	13.256	1992	1996	1997	rBV	121302	136500	1.88%	0.151%
79	13.292	1997	2002	2007	rVV2	3154746	4122991	56.69%	4.575%
80	13.341	2007	2010	2013	rVV3	181419	238023	3.27%	0.264%
81	13.372	2013	2015	2019	rVB	109058	106269	1.46%	0.118%
82	13.426	2019	2024	2031	rVB	397568	535090	7.36%	0.594%
83	13.506	2031	2037	2040	rBV2	127805	181295	2.49%	0.201%
84	13.542	2040	2043	2046	rVV	306448	386284	5.31%	0.429%
85	13.585	2046	2050	2055	rVV3	324771	540727	7.44%	0.600%
86	13.640	2055	2059	2060	rVV3	151748	181248	2.49%	0.201%
87	13.664	2060	2063	2070	rVB2	299035	478013	6.57%	0.530%
88	13.774	2074	2081	2090	rBV	2944874	3727291	51.25%	4.136%
89	13.865	2090	2096	2100	rVB2	55854	102592	1.41%	0.114%
90	13.926	2100	2106	2110	rBV2	154114	221883	3.05%	0.246%
91	13.969	2110	2113	2117	rVB	121431	138926	1.91%	0.154%
92	14.073	2126	2130	2137	rBV	221530	300293	4.13%	0.333%
93	14.213	2149	2153	2159	rBV2	202281	352794	4.85%	0.391%
94	14.323	2166	2171	2176	rBV	103882	147923	2.03%	0.164%
95	14.371	2176	2179	2184	rVB	141650	177092	2.44%	0.196%
96	14.634	2218	2222	2233	rVB	828893	1117805	15.37%	1.240%
97	14.780	2241	2246	2255	rBV	1045759	1304600	17.94%	1.448%
98	15.475	2354	2360	2366	rBV	82016	136551	1.88%	0.152%
99	15.615	2374	2383	2386	rBV	130380	210691	2.90%	0.234%
100	15.652	2386	2389	2397	rVB	73779	107520	1.48%	0.119%

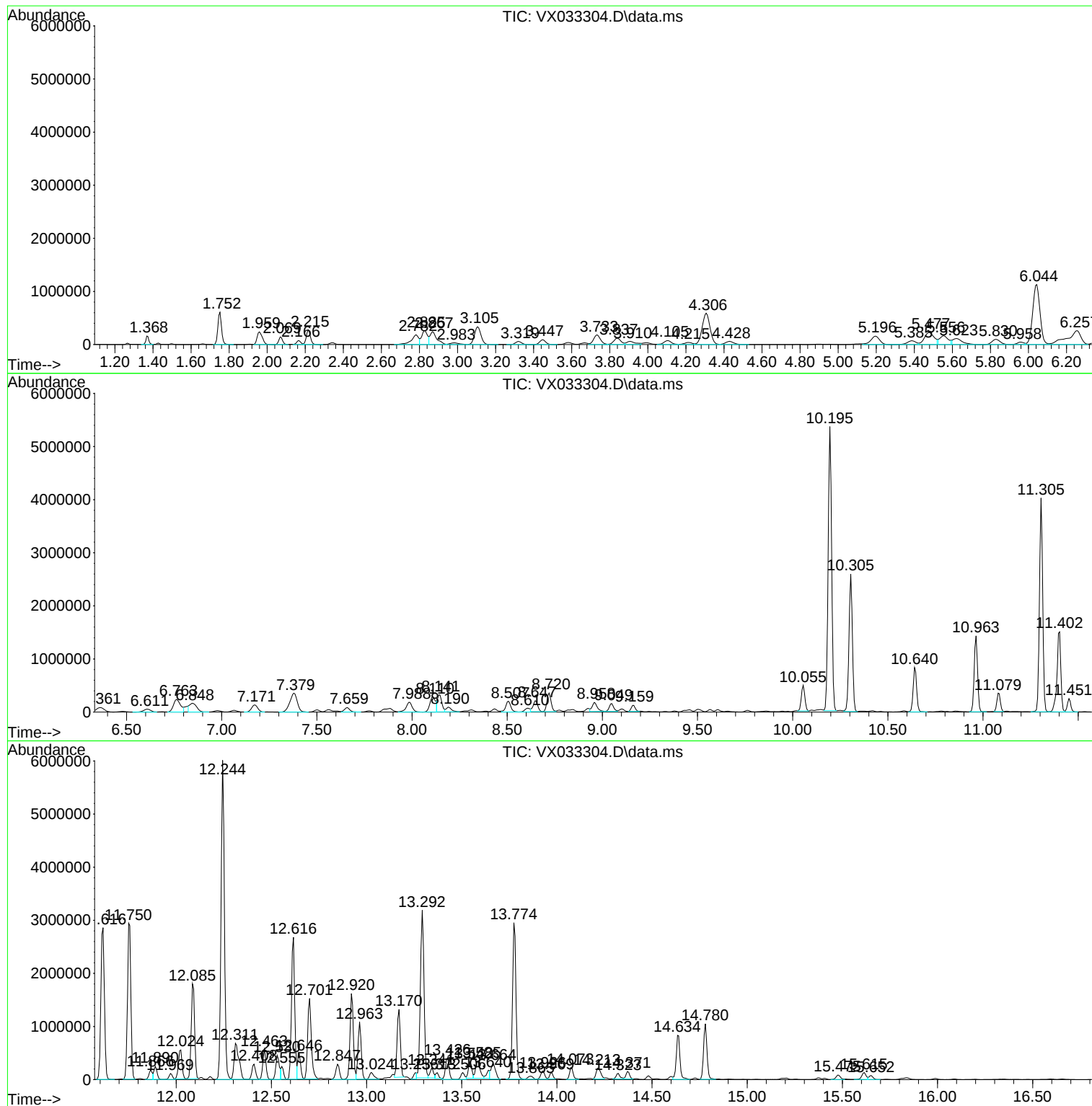
Sum of corrected areas: 90125569

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
Quant Title : SW846 8260

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



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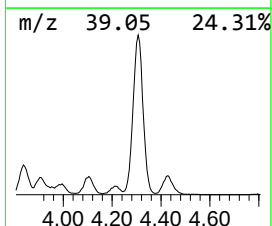
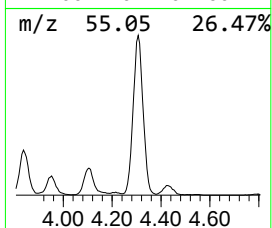
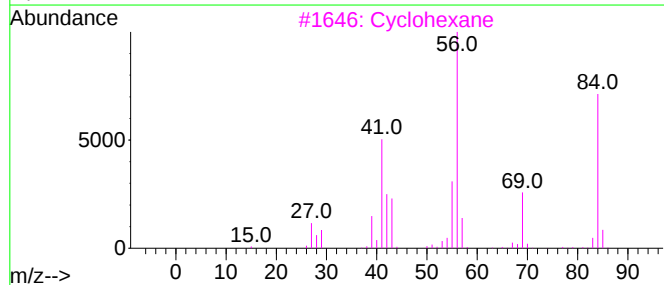
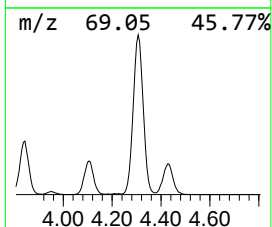
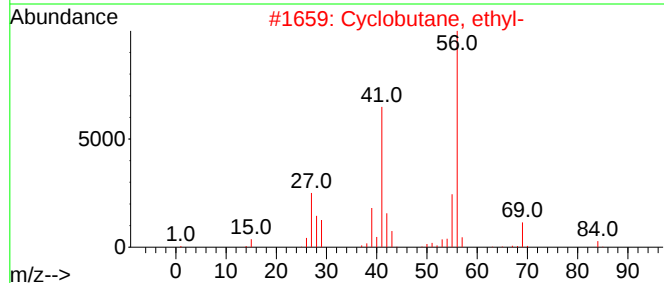
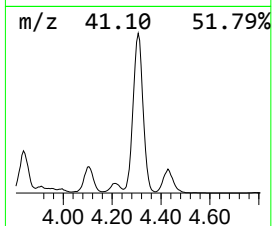
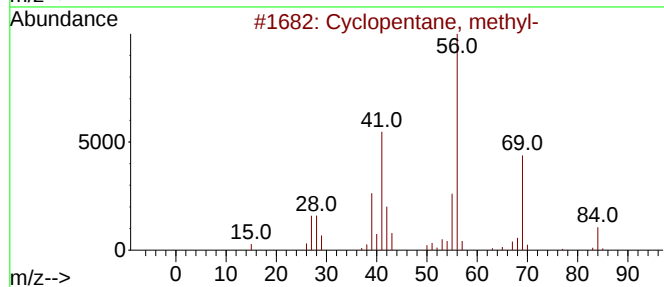
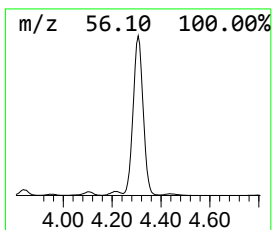
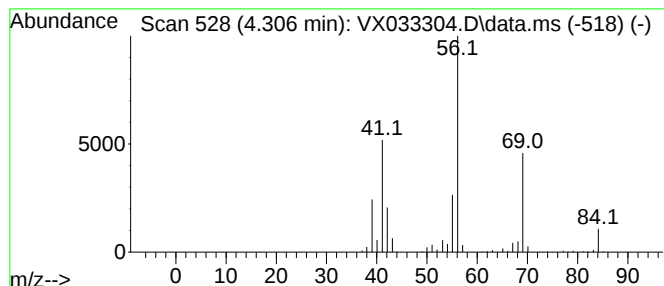
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.306	168.70 ug/l	1677190	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3			Cyclohexane	84	C6H12	000110-82-7	56
4			Cyclobutane	56	C4H8	000287-23-0	50
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43



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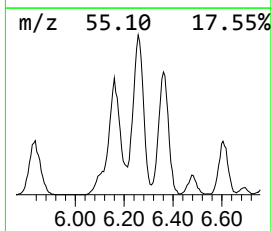
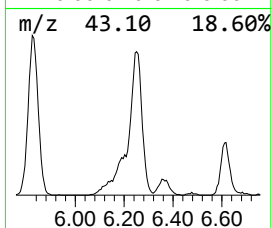
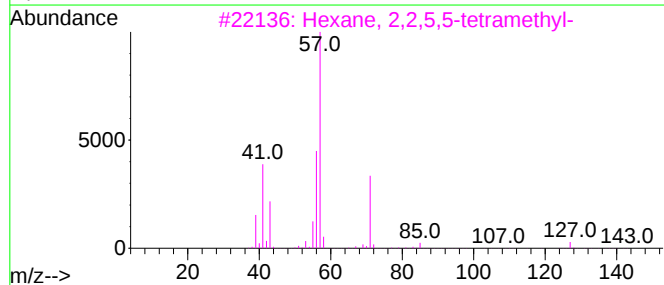
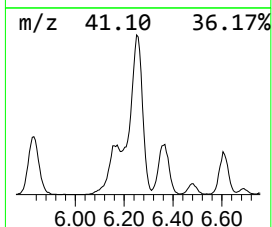
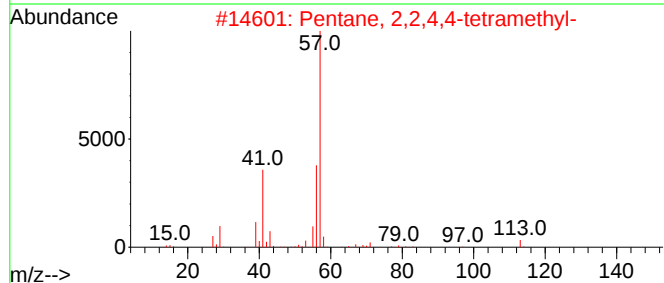
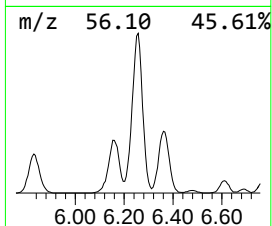
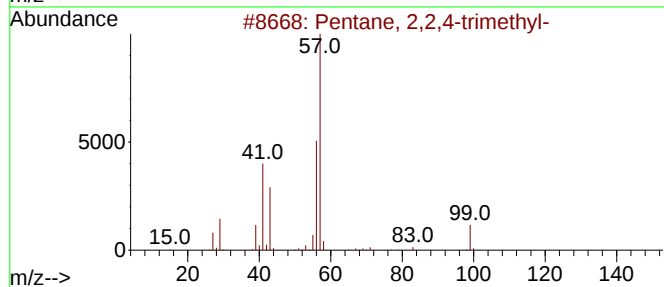
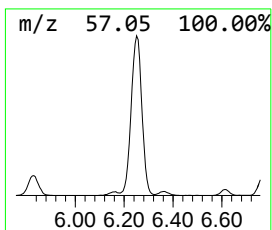
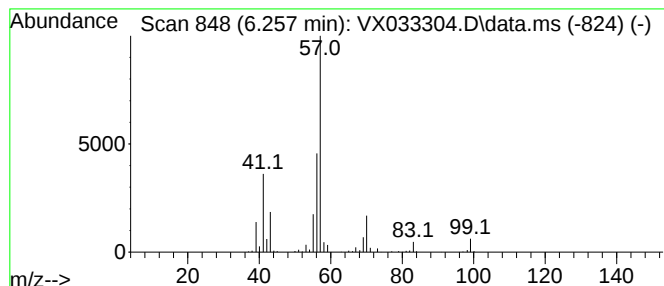
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,2,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	107.02 ug/l	1311380	1,4-Difluorobenzene	6.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	59
2			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	53
3			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	53
4			1-Butanol, 2-methyl-	88	C5H12O	000137-32-6	50
5			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	45



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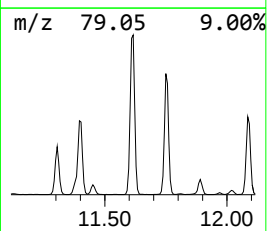
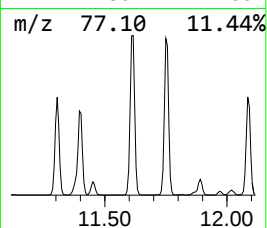
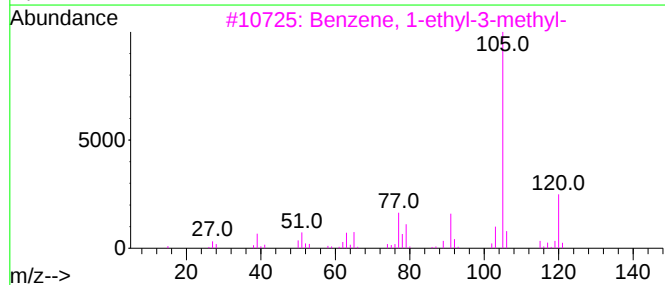
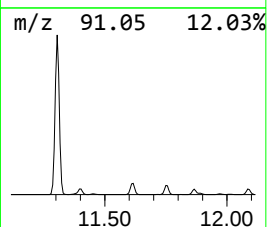
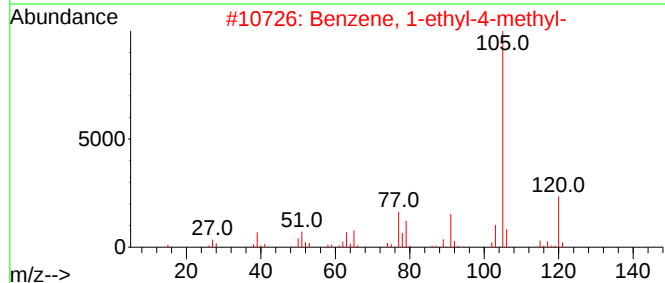
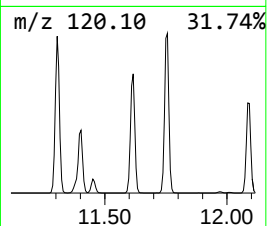
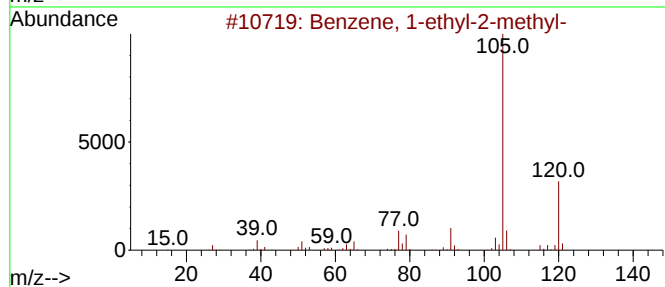
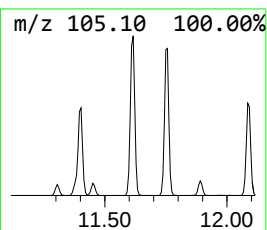
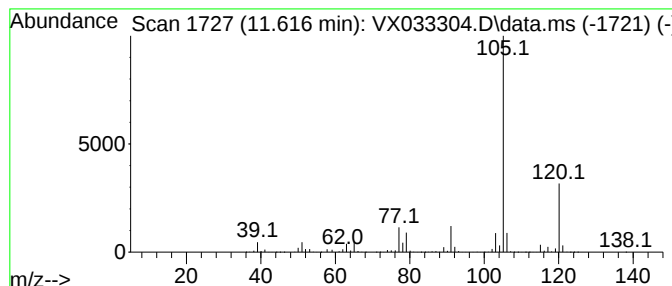
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	253.66 ug/l	3644980	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
4			Mesitylene	120	C9H12	000108-67-8	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

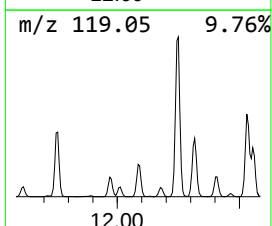
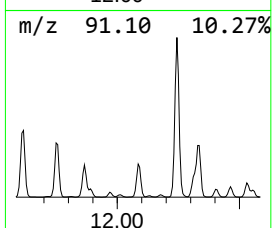
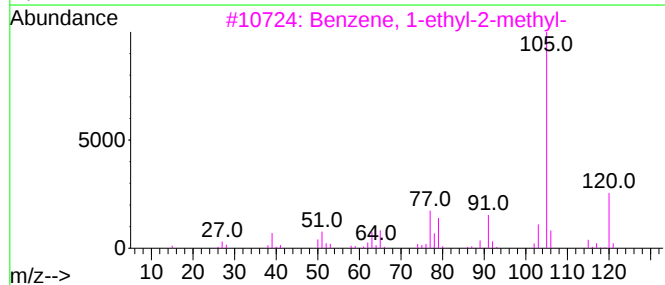
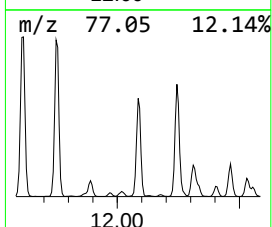
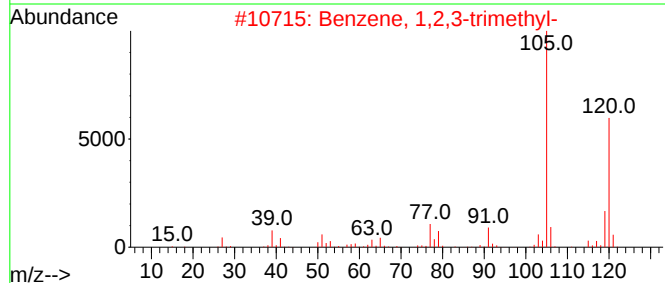
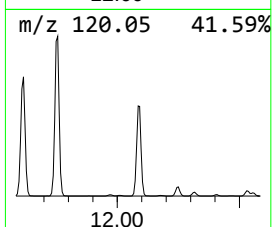
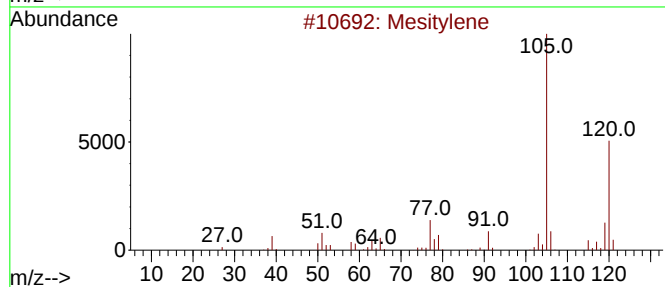
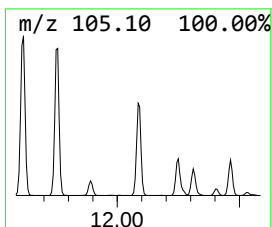
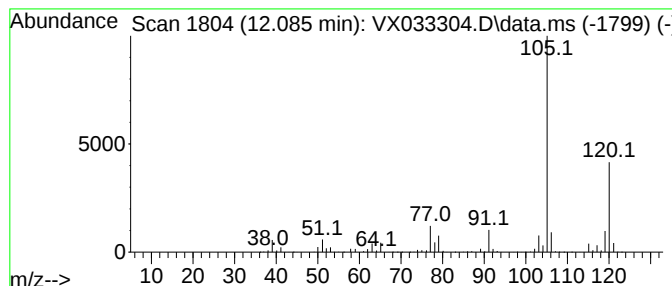
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	152.97 ug/l	2198100	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1-ethyl-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\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

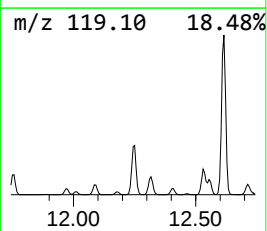
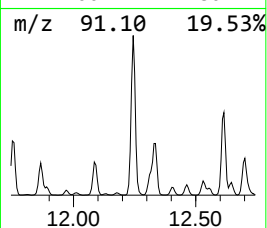
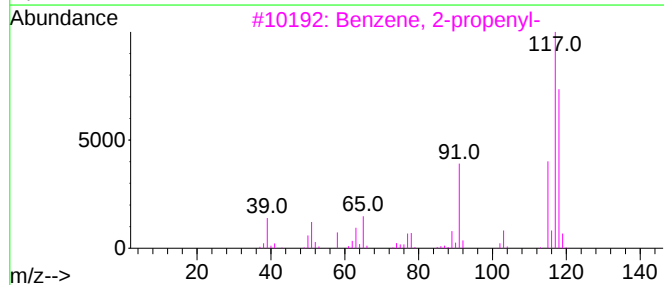
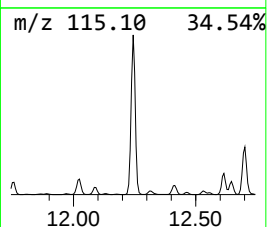
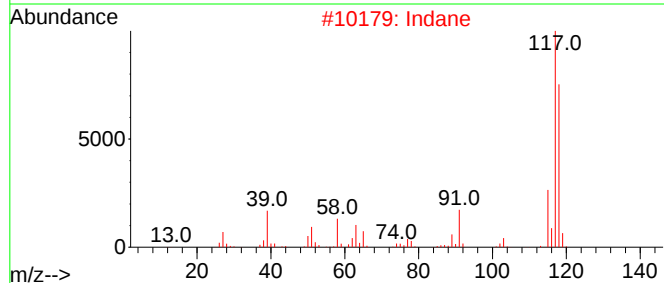
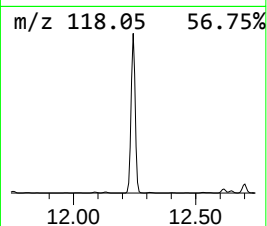
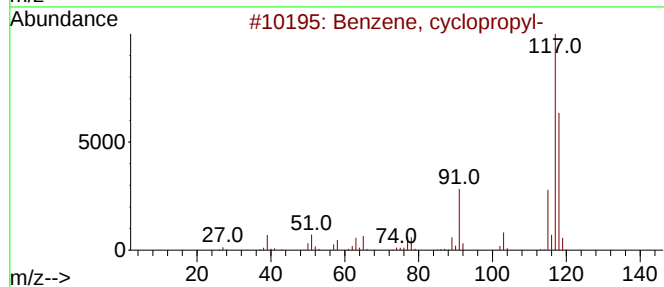
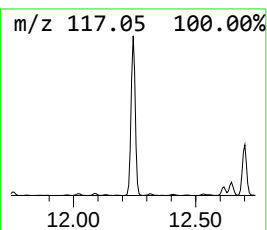
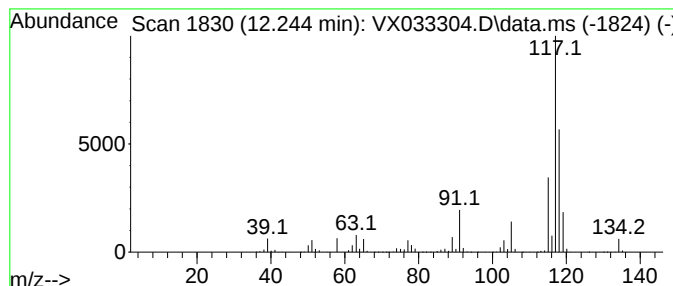
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	506.12 ug/l	7272720	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2			Indane	118	C9H10	000496-11-7	76
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4			Δ-tetralene	118	C9H10	007785-10-6	58
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

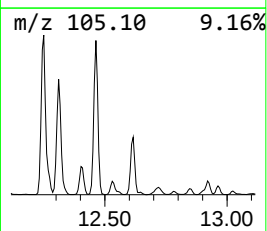
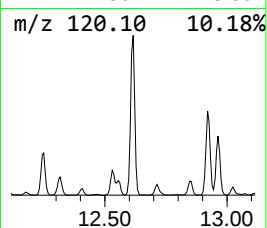
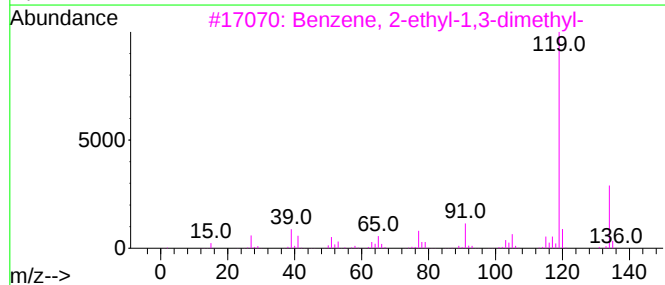
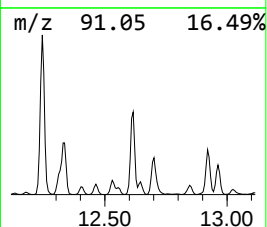
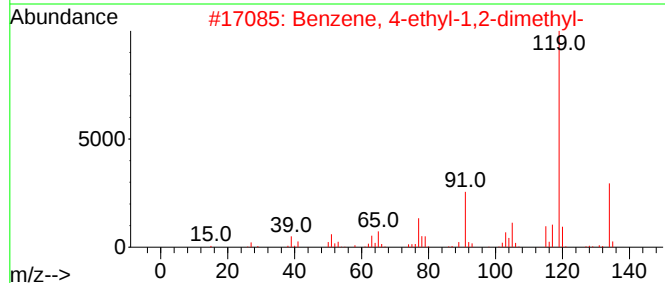
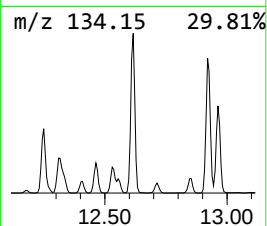
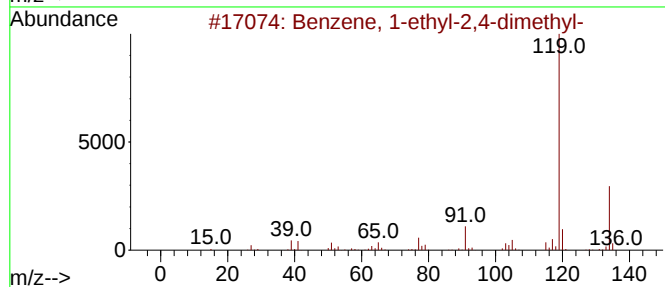
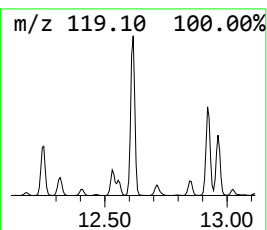
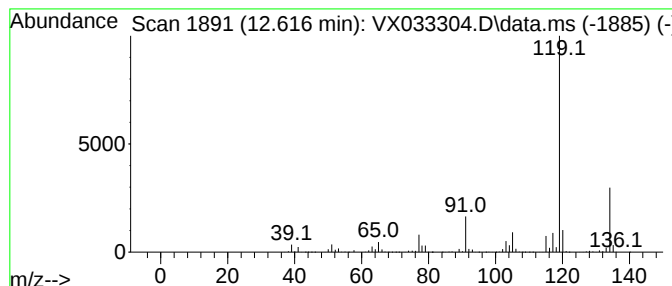
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	225.15 ug/l	3235310	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

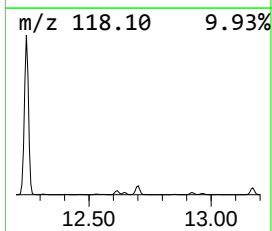
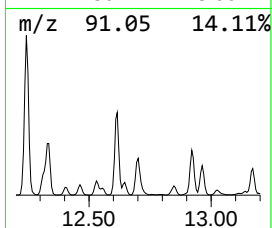
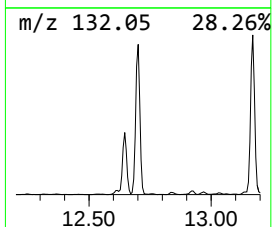
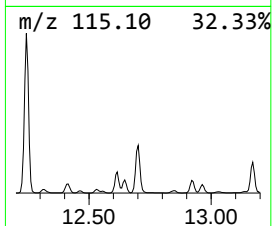
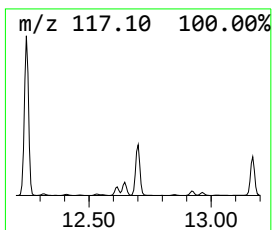
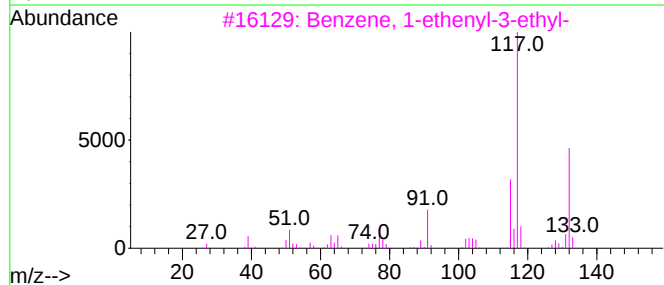
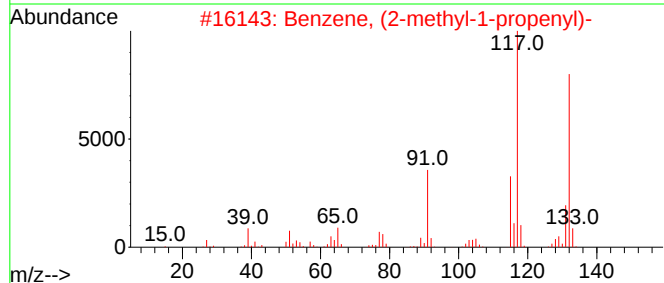
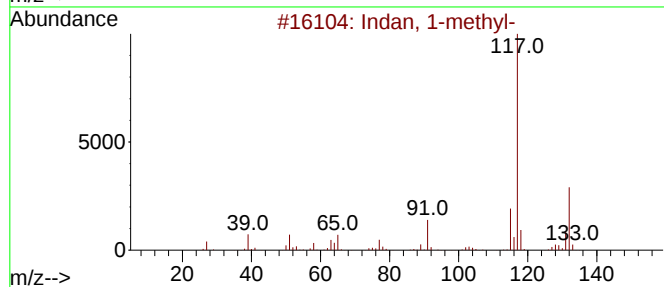
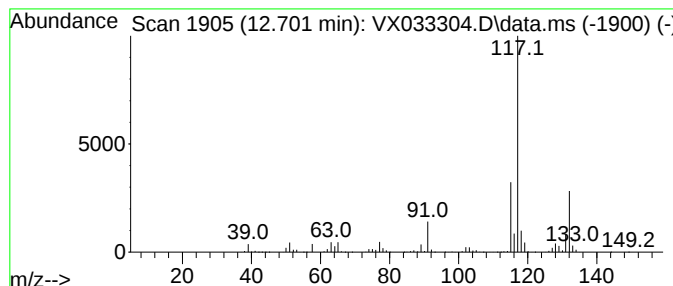
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	141.78 ug/l	2037370	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	93
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

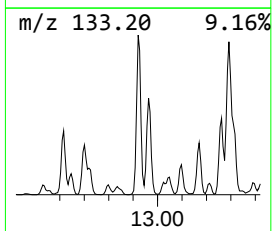
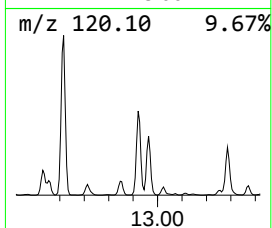
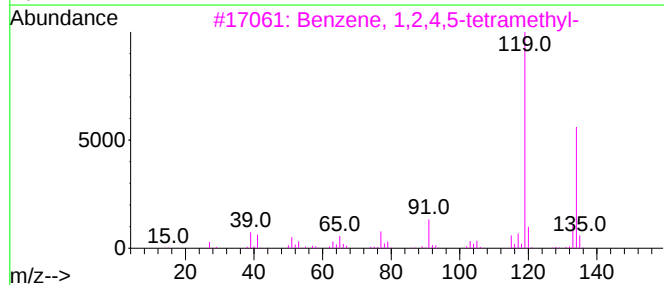
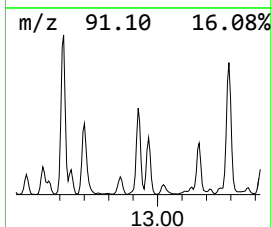
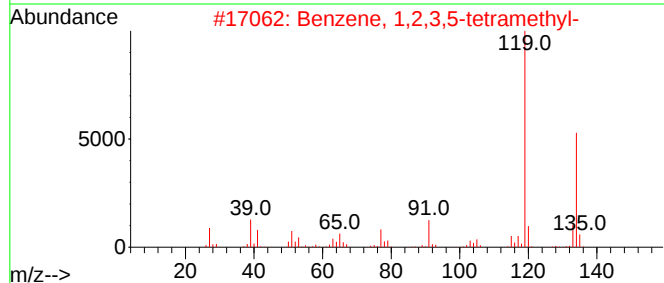
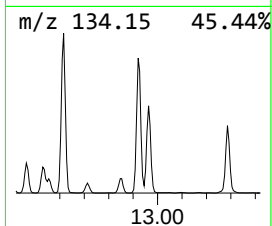
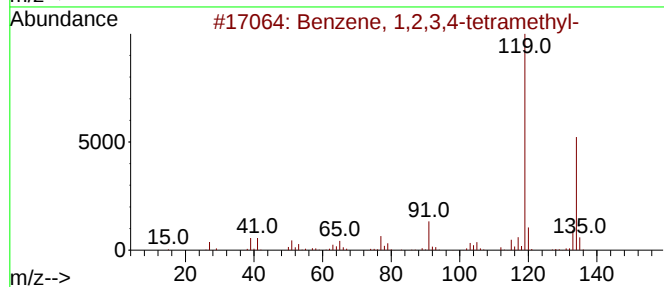
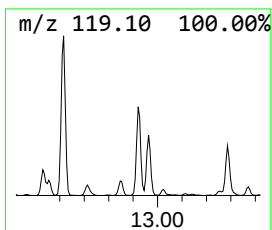
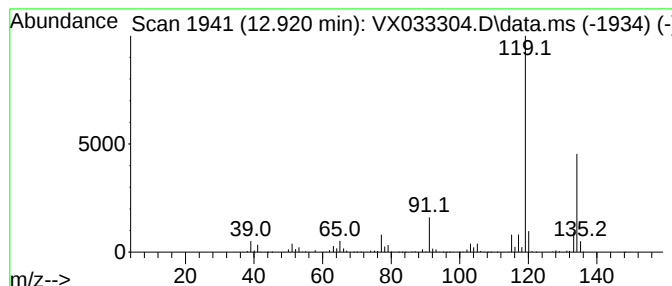
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	137.71 ug/l	1978900	1,4-Dichlorobenzene-d4	12.024

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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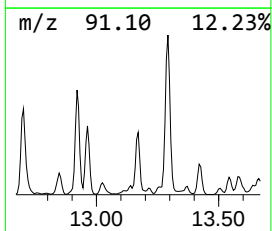
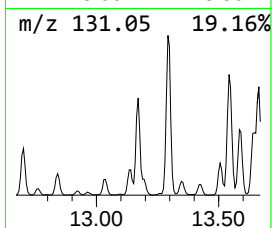
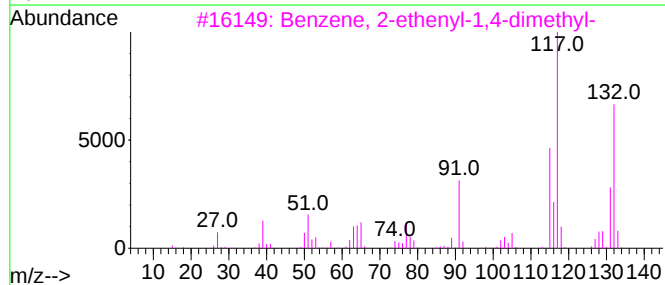
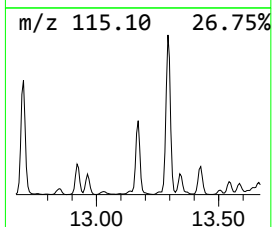
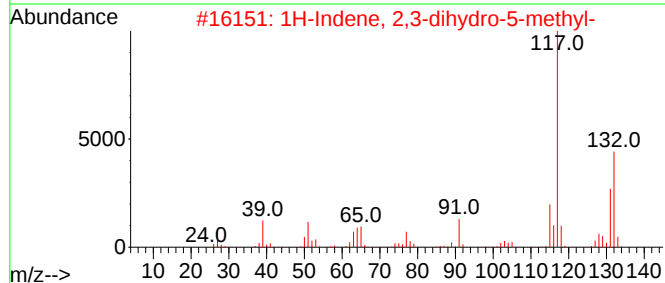
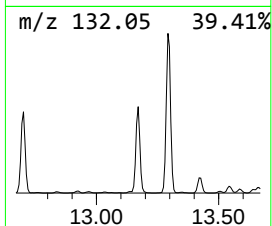
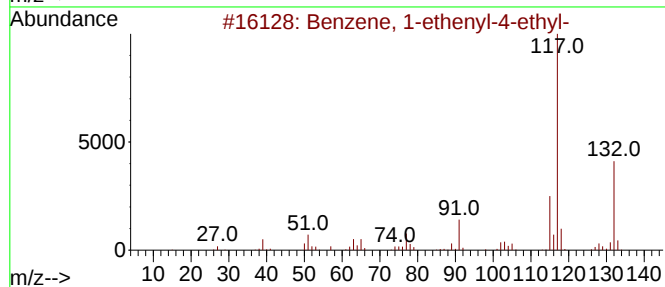
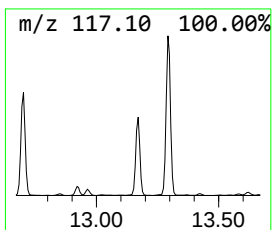
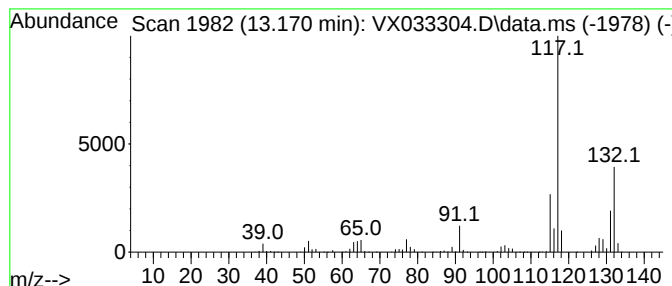
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	104.38 ug/l	1499880	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

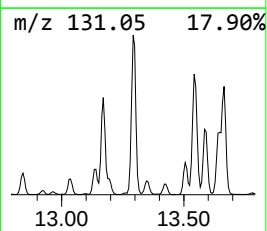
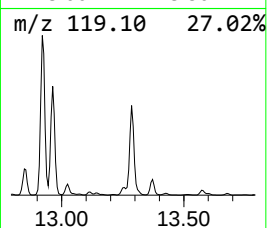
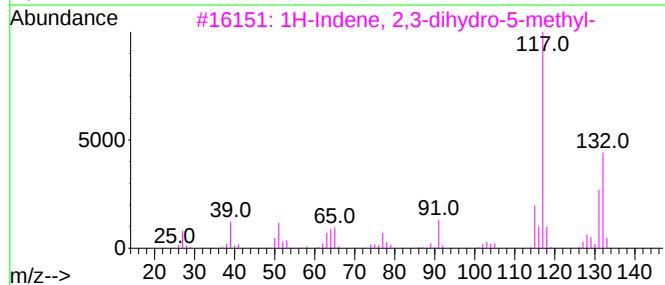
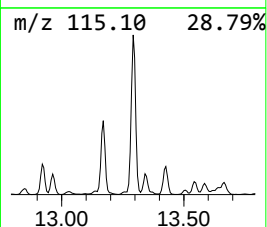
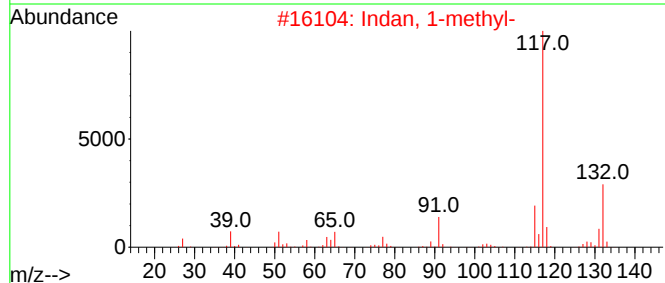
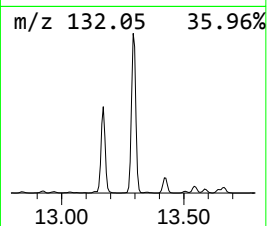
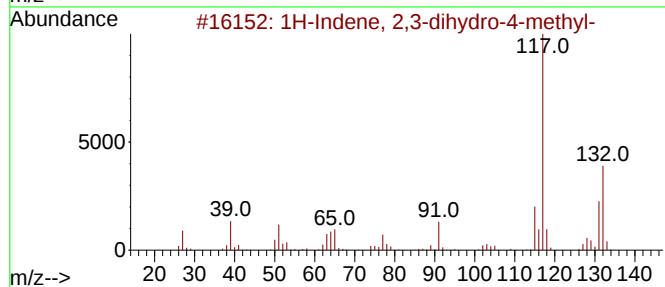
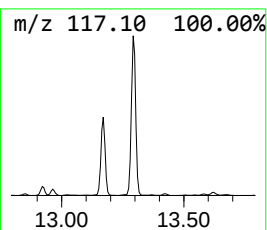
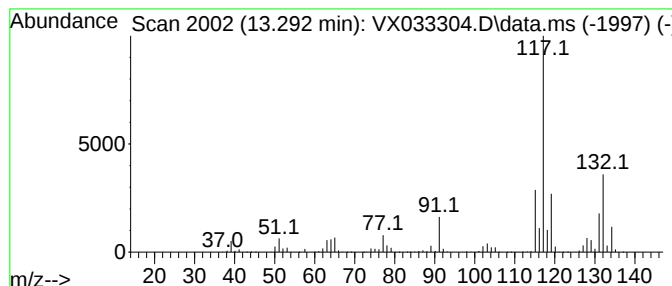
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	286.92 ug/l	4122990	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93	
2	Indan, 1-methyl-	132	C10H12	000767-58-8	93	
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89	
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81	
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120722\
Data File : VX033304.D
Acq On : 08 Dec 2022 00:10
Operator : JC/MD
Sample : N5906-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-13R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.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
Cyclopentane, m...	4.306	168.7	ug/l	1677190	1	5.556	497094	50.0
Pentane, 2,2,4-...	6.257	107.0	ug/l	1311380	2	6.763	612689	50.0
Benzene, 1-ethy...	11.616	253.7	ug/l	3644980	4	12.024	718482	50.0
Benzene, 1,2,3-...	12.085	153.0	ug/l	2198100	4	12.024	718482	50.0
Benzene, cyclop...	12.244	506.1	ug/l	7272720	4	12.024	718482	50.0
Benzene, 1-ethy...	12.616	225.2	ug/l	3235310	4	12.024	718482	50.0
Indan, 1-methyl-	12.701	141.8	ug/l	2037370	4	12.024	718482	50.0
Benzene, 1,2,3,...	12.920	137.7	ug/l	1978900	4	12.024	718482	50.0
Benzene, 1-ethe...	13.170	104.4	ug/l	1499880	4	12.024	718482	50.0
1H-Indene, 2,3-...	13.292	286.9	ug/l	4122990	4	12.024	718482	50.0