

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
 Data File : VX031549.D
 Acq On : 22 Sep 2022 04:56
 Operator : JC/MD
 Sample : N4642-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-3R

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\82X091922W.M
 Title : SW846 8260

Signal : TIC: VX031549.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.337	198	205	211	rBV2	6098	11318	1.73%	0.207%
2	2.782	270	278	283	rBV3	6881	14230	2.17%	0.260%
3	3.099	322	330	339	rBV3	4409	9884	1.51%	0.181%
4	5.391	693	706	719	rBV	77318	220134	33.60%	4.028%
5	5.556	723	733	742	rBV2	74652	209314	31.95%	3.830%
6	5.842	771	780	789	rVB6	2665	7062	1.08%	0.129%
7	5.958	789	799	806	rVV2	82688	218375	33.34%	3.996%
8	6.044	806	813	823	rVV	47339	125849	19.21%	2.303%
9	6.159	823	832	841	rVV6	6623	19997	3.05%	0.366%
10	6.257	841	848	858	rVV3	5848	15406	2.35%	0.282%
11	6.763	922	931	942	rBV2	116988	273858	41.80%	5.011%
12	7.373	1021	1031	1041	rBV6	6388	17193	2.62%	0.315%
13	7.958	1121	1127	1135	rVB5	3344	7276	1.11%	0.133%
14	8.324	1178	1187	1192	rBV6	3816	7654	1.17%	0.140%
15	8.604	1227	1233	1235	rBV4	9684	16997	2.59%	0.311%
16	8.647	1235	1240	1253	rVB	412960	655086	100.00%	11.987%
17	8.775	1256	1261	1266	rBV3	8426	14555	2.22%	0.266%
18	8.958	1286	1291	1300	rVB5	32837	72590	11.08%	1.328%
19	9.049	1300	1306	1311	rVV	34073	54226	8.28%	0.992%
20	9.104	1311	1315	1320	rVB2	13549	21649	3.30%	0.396%
21	9.159	1320	1324	1336	rVB	12904	22101	3.37%	0.404%
22	9.458	1364	1373	1377	rBV5	3921	7147	1.09%	0.131%
23	9.513	1377	1382	1388	rVV4	10243	17287	2.64%	0.316%
24	10.055	1465	1471	1477	rBV	260567	344007	52.51%	6.295%
25	10.128	1480	1483	1491	rVV	8717	13885	2.12%	0.254%
26	10.195	1491	1494	1499	rVV2	5831	8178	1.25%	0.150%
27	10.275	1502	1507	1510	rVV5	6161	11348	1.73%	0.208%
28	10.457	1533	1537	1542	rVB4	4586	7169	1.09%	0.131%
29	10.604	1551	1561	1565	rBV3	7086	16790	2.56%	0.307%
30	10.964	1616	1620	1627	rBV2	6136	9714	1.48%	0.178%
31	11.079	1634	1639	1645	rBV	351439	440984	67.32%	8.069%
32	11.616	1721	1727	1734	rVB	13432	17750	2.71%	0.325%
33	11.866	1762	1768	1769	rBV	7749	9058	1.38%	0.166%
34	11.890	1769	1772	1778	rVV2	14314	19131	2.92%	0.350%
35	11.969	1781	1785	1789	rVV3	6670	8747	1.34%	0.160%
36	12.024	1789	1794	1802	rVV	265119	332499	50.76%	6.084%

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37	12.177	1815	1819	1823	rVB	5294	6564	1.00%	0.120%
38	12.244	1825	1830	1837	rVV2	164019	212912	32.50%	3.896%
39	12.317	1837	1842	1849	rVB	29815	40825	6.23%	0.747%
40	12.408	1851	1857	1862	rBV2	14334	19655	3.00%	0.360%
41	12.530	1871	1877	1885	rBV3	3950	8103	1.24%	0.148%
42	12.616	1885	1891	1893	rVV	26947	34862	5.32%	0.638%
43	12.646	1893	1896	1900	rVV	46261	57258	8.74%	1.048%
44	12.701	1900	1905	1912	rVV	224465	281852	43.03%	5.157%
45	12.762	1912	1915	1918	rVV2	7685	10275	1.57%	0.188%
46	12.841	1921	1928	1933	rVB	15306	21762	3.32%	0.398%
47	12.920	1938	1941	1946	rVB2	15029	19609	2.99%	0.359%
48	13.030	1954	1959	1965	rBV3	32118	49558	7.57%	0.907%
49	13.134	1965	1976	1979	rBV3	35204	61743	9.43%	1.130%
50	13.170	1979	1982	1985	rBV	47522	48918	7.47%	0.895%
51	13.250	1992	1995	1998	rBV2	13436	18803	2.87%	0.344%
52	13.292	1998	2002	2008	rVV	250070	329215	50.26%	6.024%
53	13.347	2008	2011	2012	rVV2	18667	23034	3.52%	0.421%
54	13.372	2012	2015	2020	rVV	34709	47407	7.24%	0.867%
55	13.426	2020	2024	2030	rVB	30358	40422	6.17%	0.740%
56	13.506	2030	2037	2040	rBV2	22308	33191	5.07%	0.607%
57	13.542	2040	2043	2047	rVV	85327	118228	18.05%	2.163%
58	13.585	2047	2050	2054	rVV2	82703	117245	17.90%	2.145%
59	13.640	2054	2059	2061	rVV3	58720	108591	16.58%	1.987%
60	13.664	2061	2063	2072	rVB2	91188	106017	16.18%	1.940%
61	13.756	2072	2078	2079	rBV2	7856	10616	1.62%	0.194%
62	13.798	2079	2085	2091	rVV4	10106	29012	4.43%	0.531%
63	13.859	2091	2095	2099	rVV4	11041	18649	2.85%	0.341%
64	13.926	2099	2106	2110	rVV2	29747	43178	6.59%	0.790%
65	13.969	2110	2113	2117	rVV	36118	46824	7.15%	0.857%
66	14.006	2117	2119	2126	rVB2	9519	12281	1.87%	0.225%
67	14.079	2126	2131	2135	rBV	31790	41396	6.32%	0.757%
68	14.213	2149	2153	2160	rVV3	26590	47156	7.20%	0.863%
69	14.323	2168	2171	2175	rVV3	9956	10744	1.64%	0.197%
70	14.371	2175	2179	2183	rVB	34716	42433	6.48%	0.776%
71	14.420	2183	2187	2193	rVB2	6395	8503	1.30%	0.156%
72	14.487	2193	2198	2202	rBV3	13174	19743	3.01%	0.361%
73	14.603	2210	2217	2221	rBV2	10099	20450	3.12%	0.374%
74	14.780	2242	2246	2251	rBV4	12434	19596	2.99%	0.359%

Sum of corrected areas: 5465078

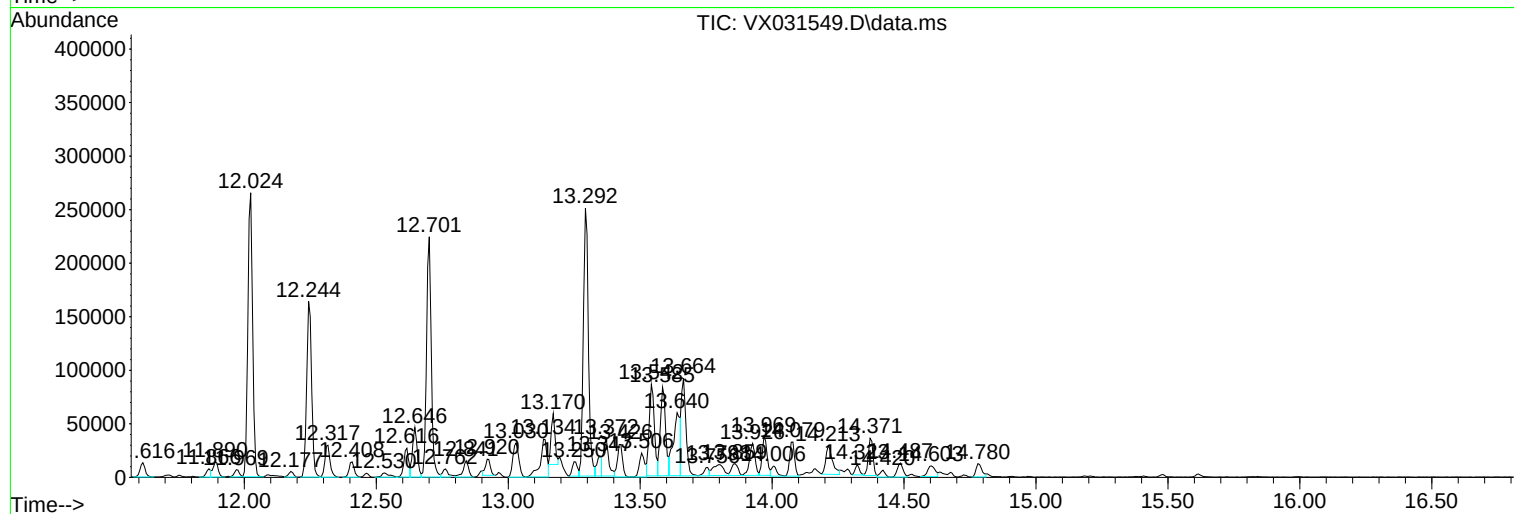
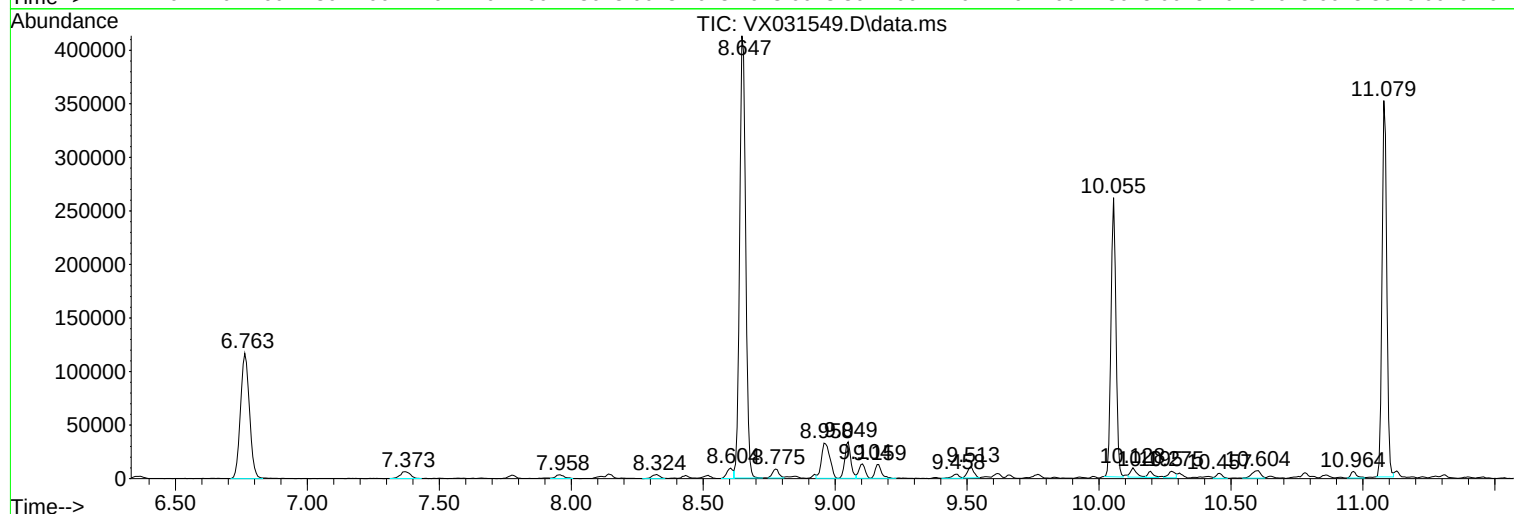
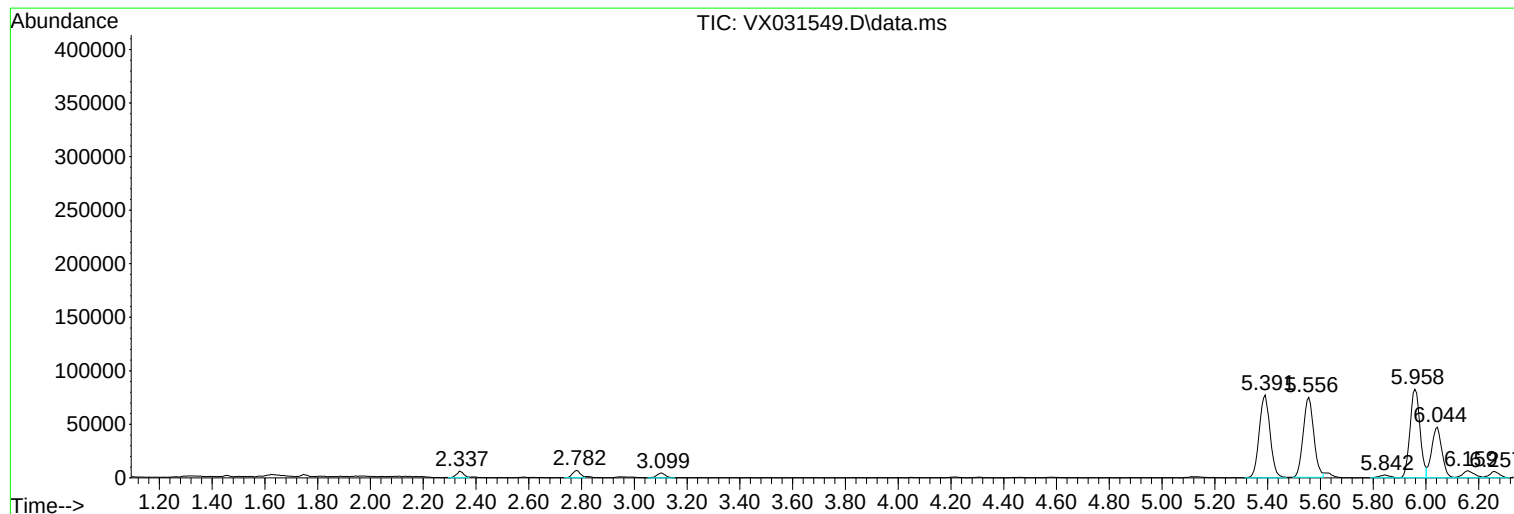
Instrument :
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Instrument :
MSVOA_X
ClientSampleId :
MW-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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



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Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3R

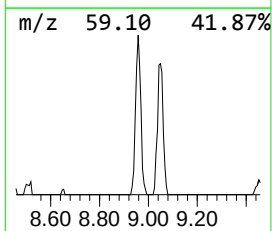
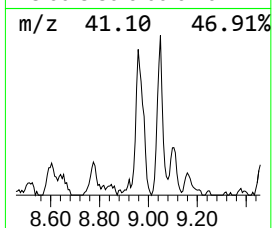
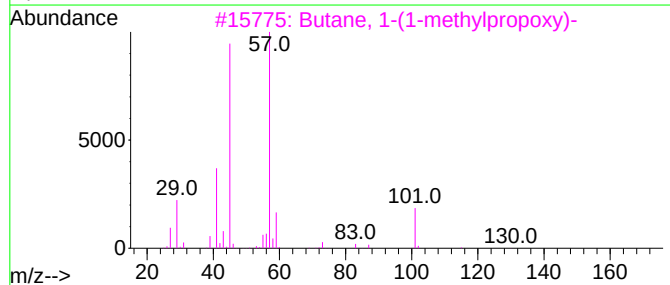
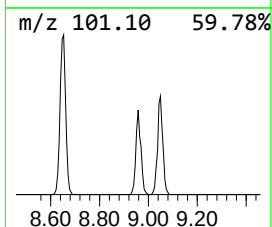
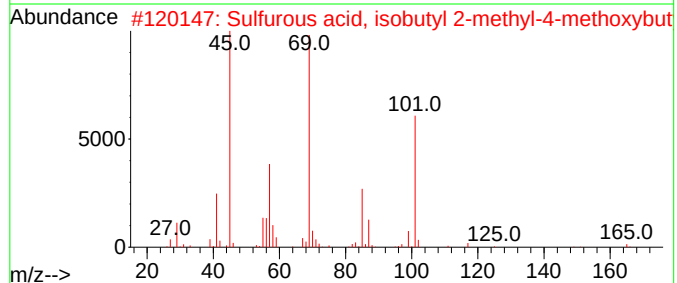
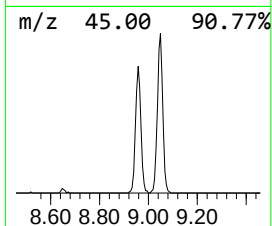
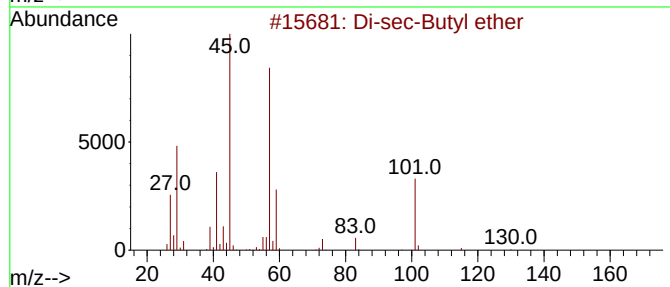
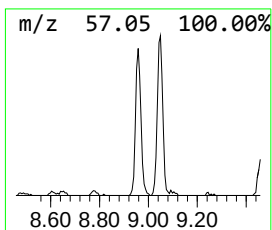
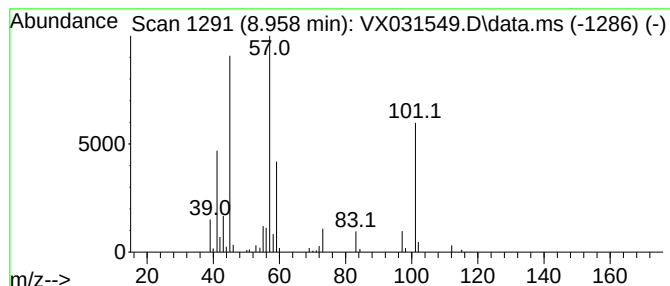
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 1 Di-sec-Butyl ether Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.958	10.55 ug/l	72590	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Di-sec-Butyl ether	130	C8H18O	006863-58-7	83
2			Sulfurous acid, isobutyl 2-methy...	238	C10H22O4S	1000309-20-3	50
3			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	45
4			2-Octanol, (S)-	130	C8H18O	006169-06-8	43
5			2-Octanol	130	C8H18O	000123-96-6	43



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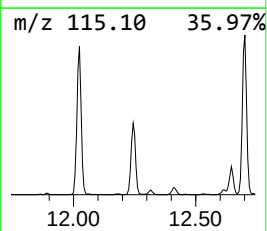
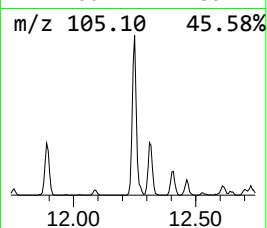
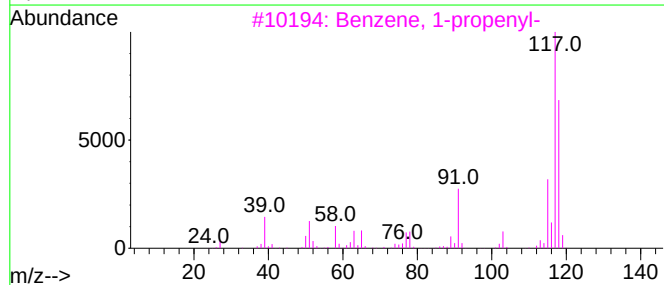
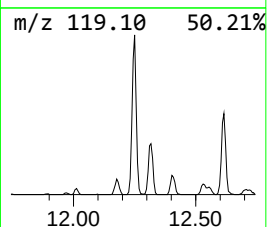
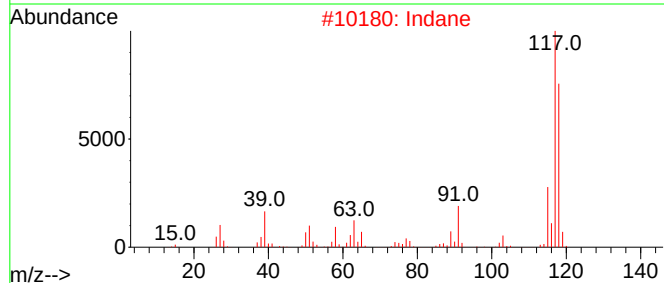
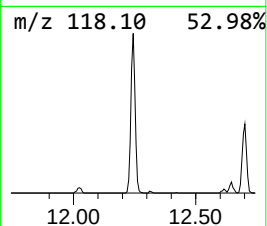
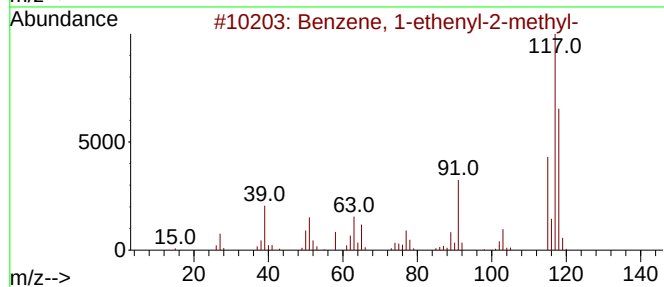
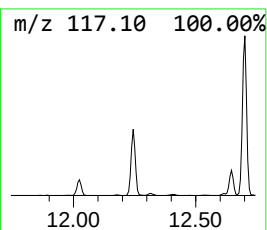
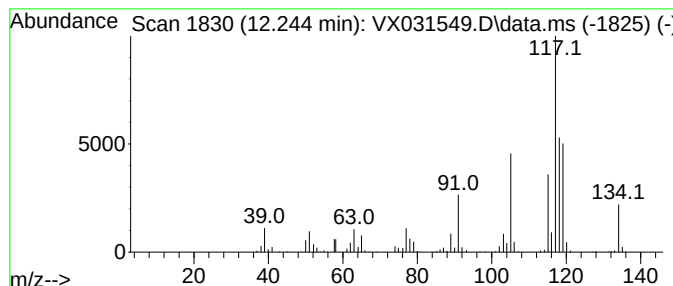
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	32.02 ug/l	212912	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
2			Indane	118	C9H10	000496-11-7	60
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	55
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	55



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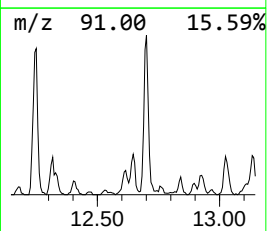
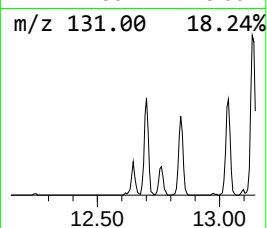
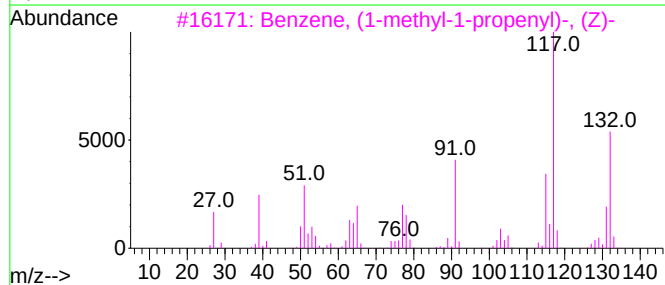
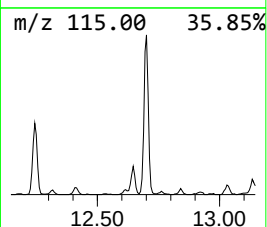
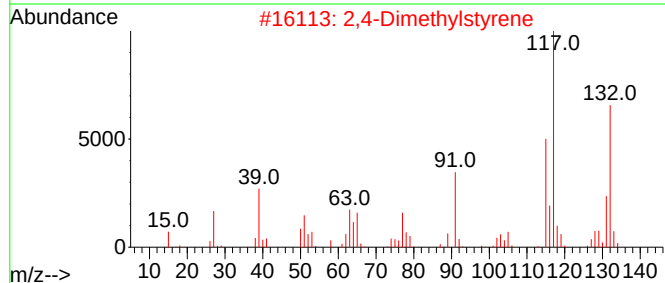
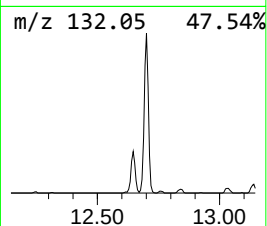
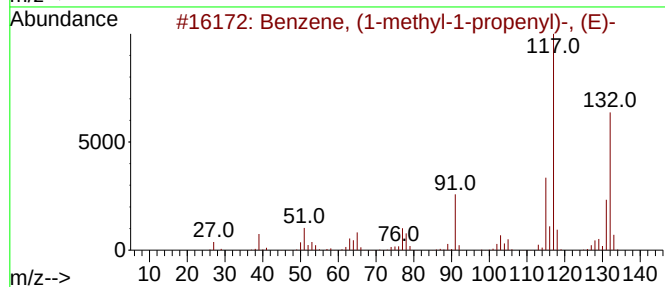
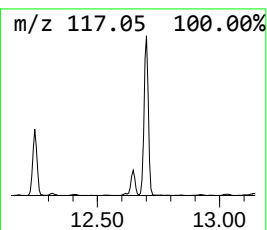
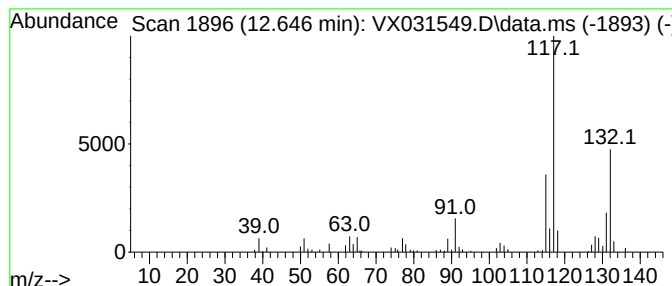
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, (1-methyl-1-propenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	8.61 ug/l	57258	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	94
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	93
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	91



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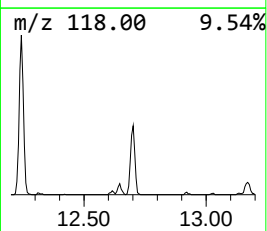
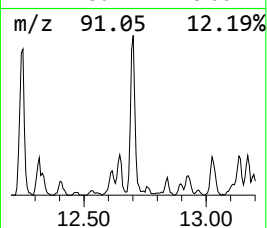
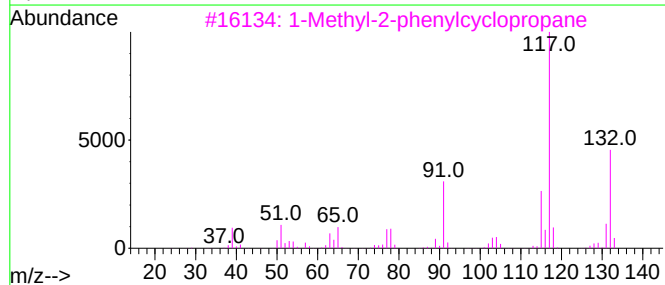
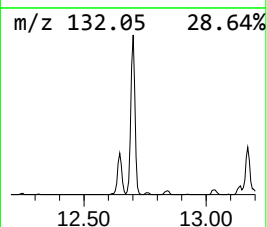
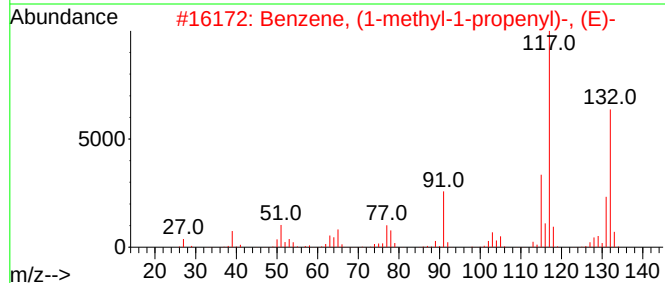
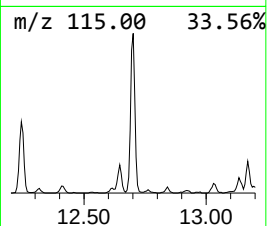
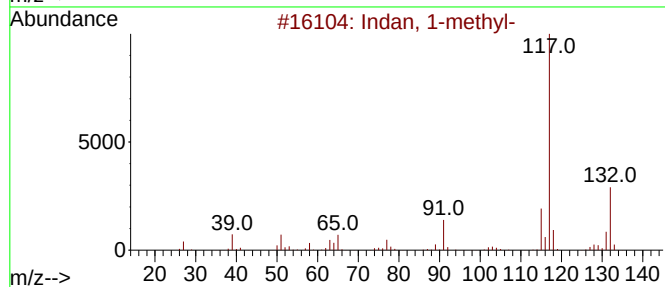
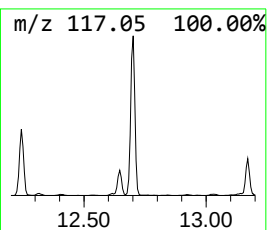
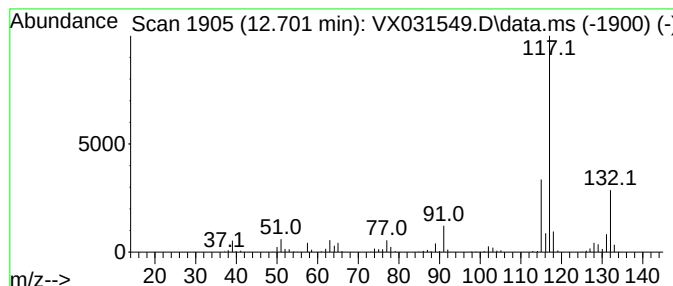
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Quant Title : SW846 8260

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

Peak Number 4 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	42.38 ug/l	281852	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	90
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3R

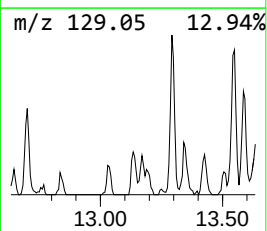
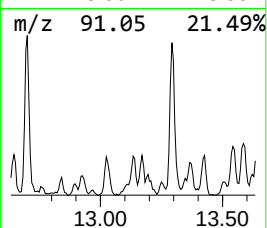
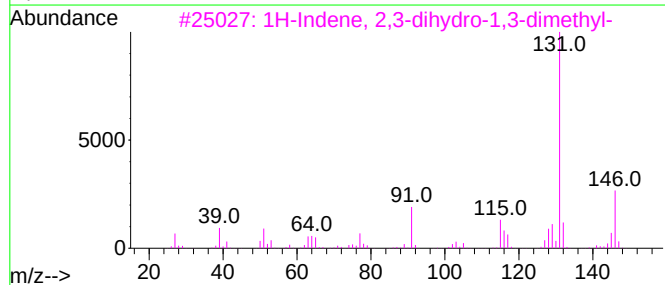
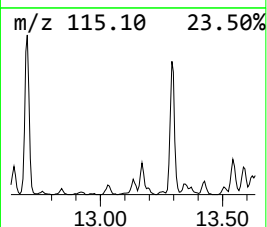
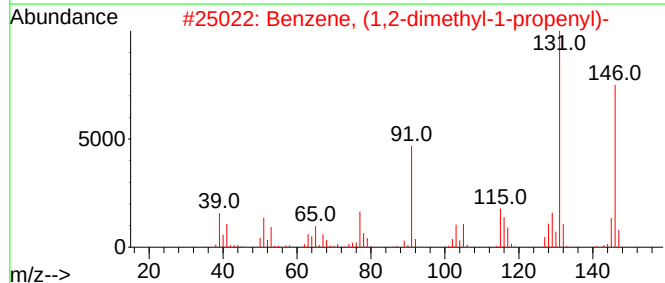
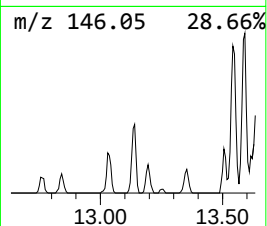
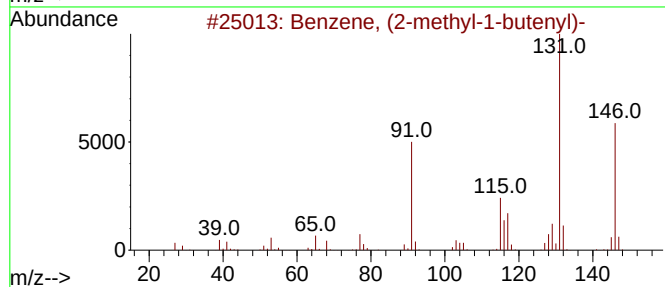
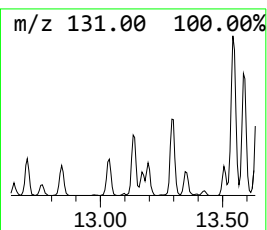
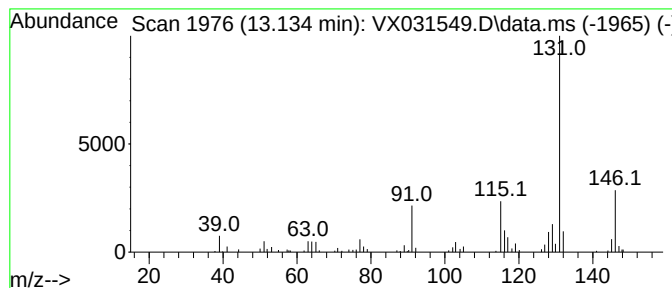
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.134	9.28 ug/l	61743	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3R

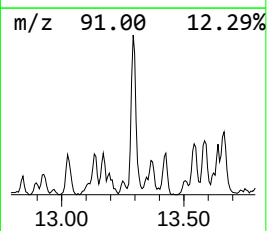
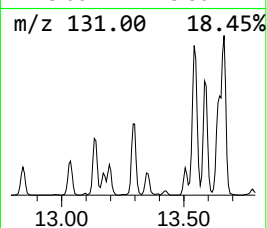
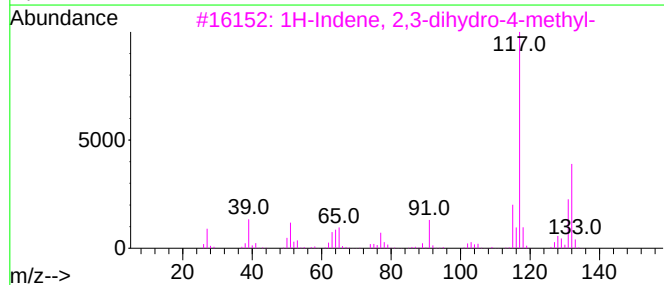
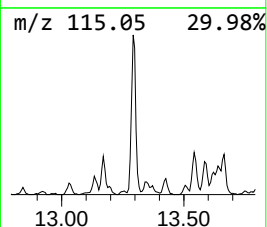
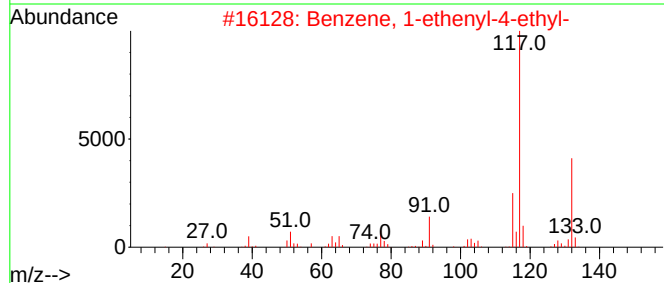
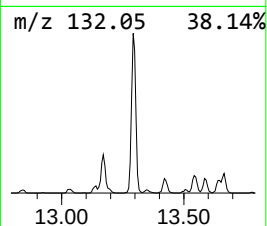
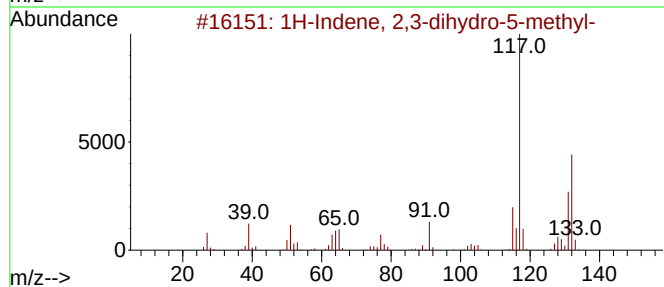
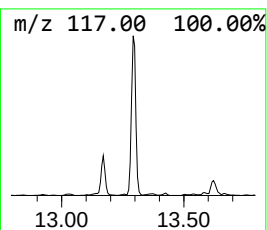
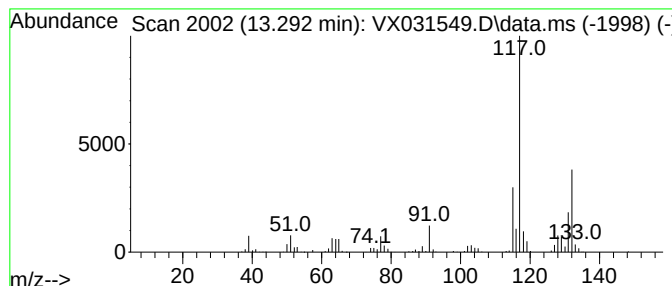
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	49.51 ug/l	329215	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93	
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90	
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90	
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90	
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3R

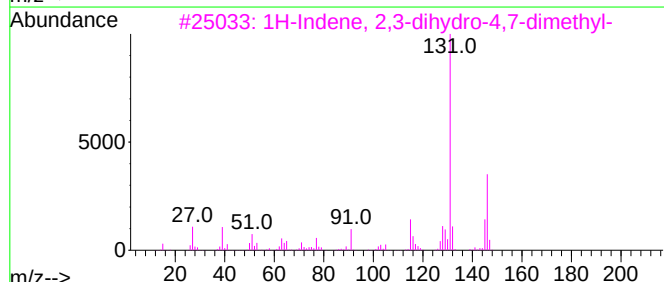
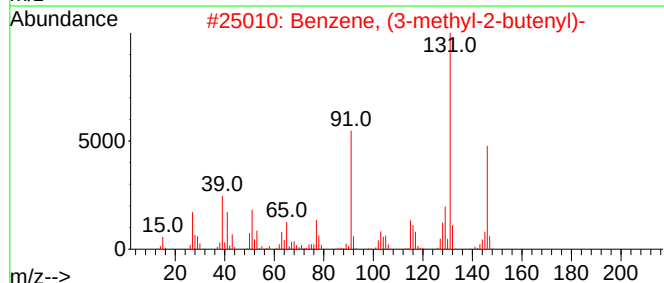
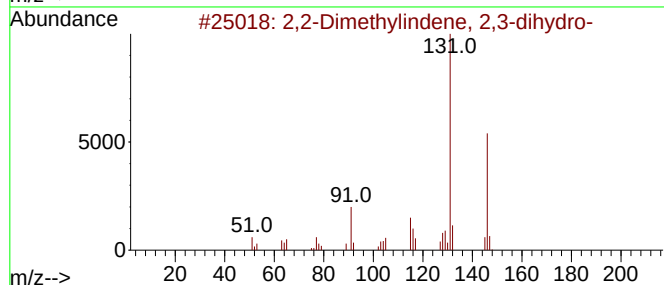
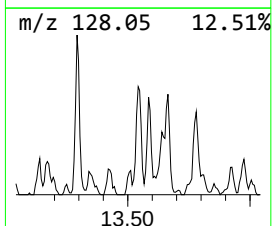
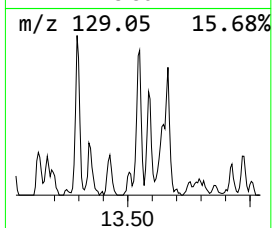
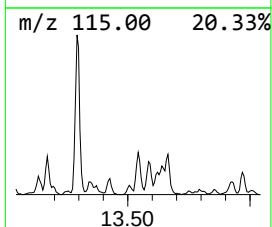
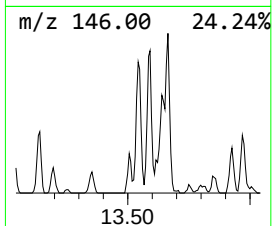
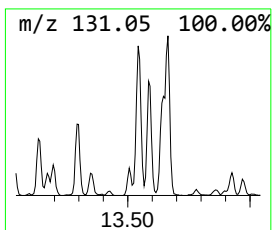
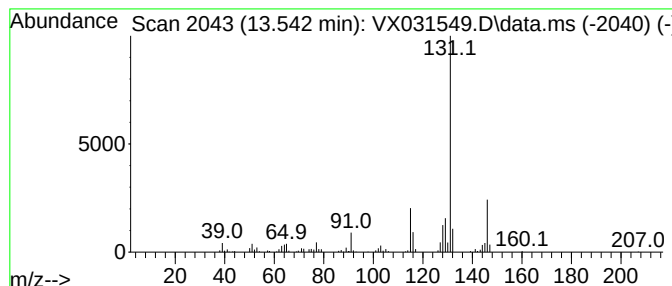
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Quant Title : SW846 8260

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

Peak Number 7 2,2-Dimethylindene, 2,3-dih... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	17.78 ug/l	118228	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	91
5		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

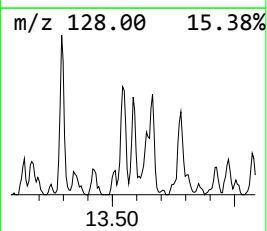
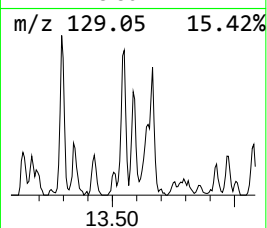
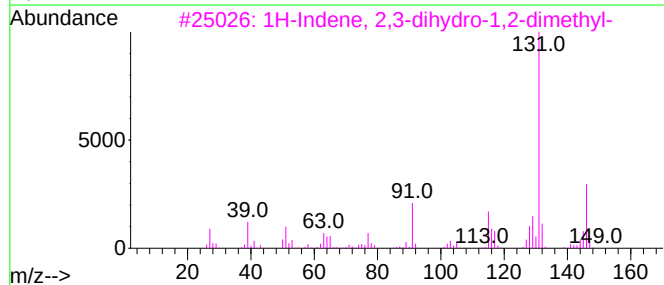
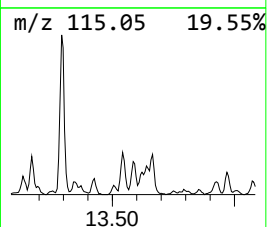
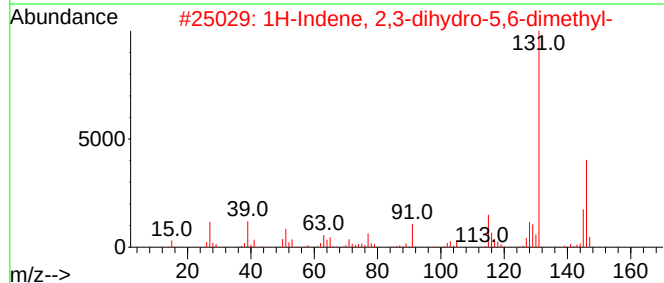
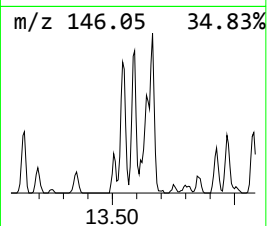
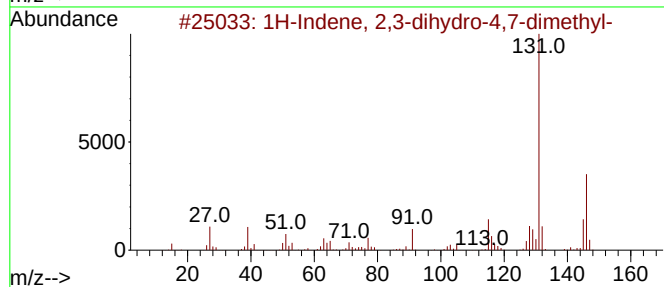
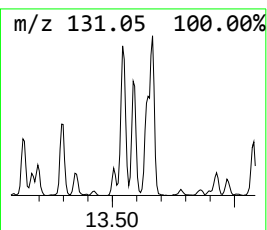
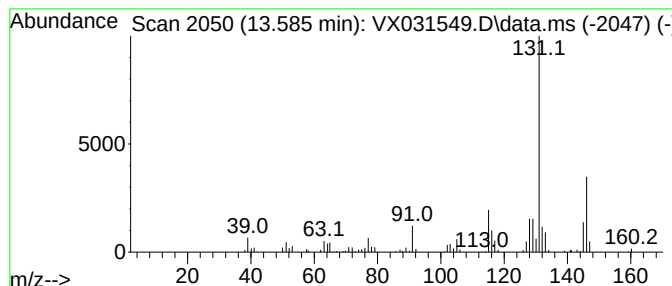
Instrument :
MSVOA_X
ClientSampleId :
MW-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.585	17.63 ug/l	117245	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...		146	C11H14	006682-71-9	95
2	1H-Indene, 2,3-dihydro-5,6-dimet...		146	C11H14	001075-22-5	93
3	1H-Indene, 2,3-dihydro-1,2-dimet...		146	C11H14	017057-82-8	93
4	Benzene, (1-methyl-1-butenyl)-		146	C11H14	053172-84-2	91
5	2,2-Dimethylindene, 2,3-dihydro-		146	C11H14	020836-11-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

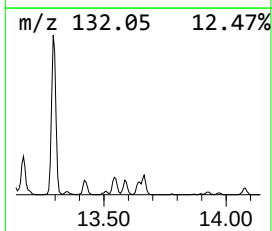
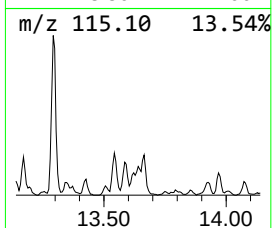
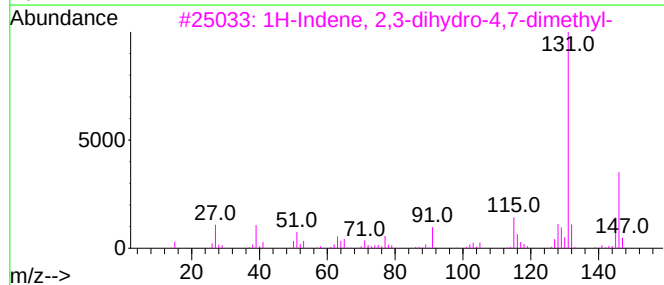
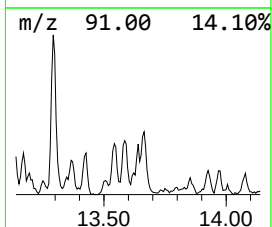
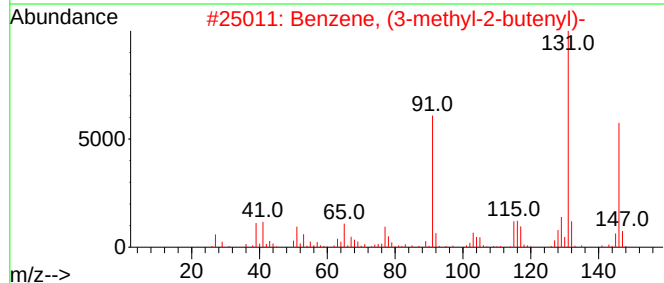
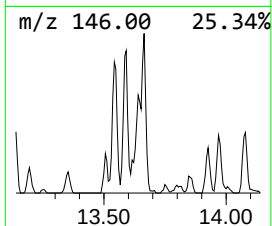
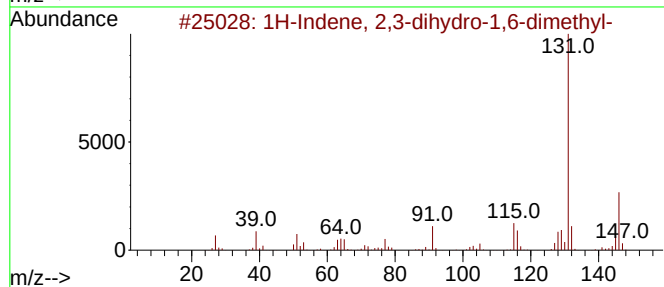
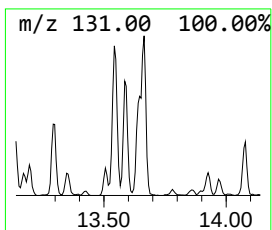
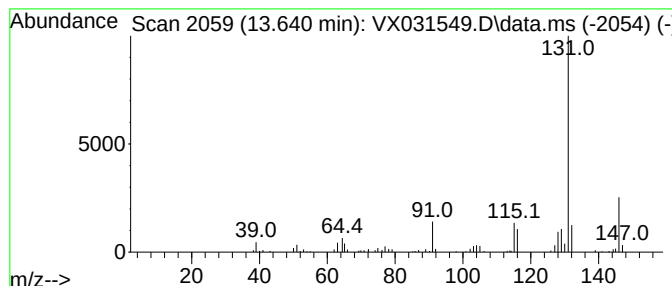
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.640	16.33 ug/l	108591	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...		146	C11H14	017059-48-2	94
2	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	91
3	1H-Indene, 2,3-dihydro-4,7-dimet...		146	C11H14	006682-71-9	91
4	Benzene, 1-methyl-2-(1-methyl-2-...		146	C11H14	097664-19-2	91
5	Benzene, 1-methyl-4-(1-methyl-2-...		146	C11H14	097664-18-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-3R

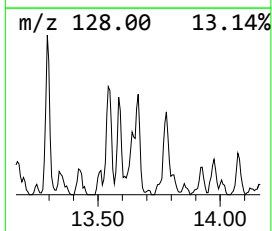
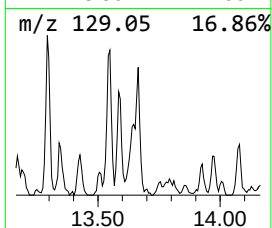
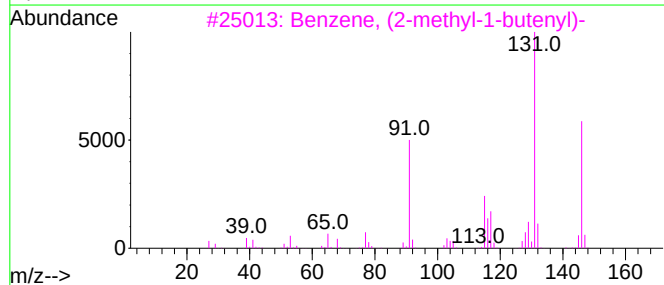
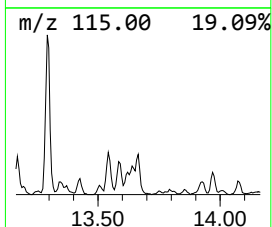
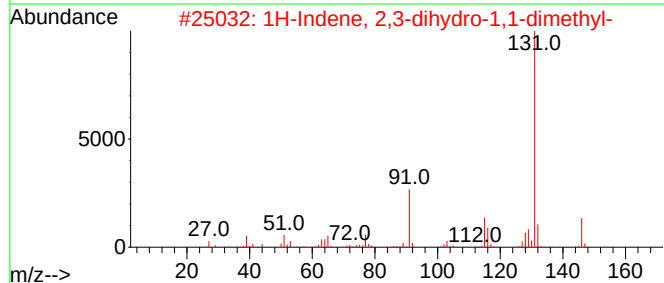
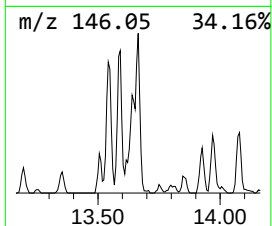
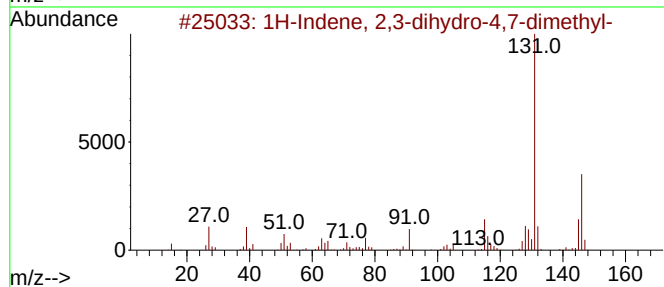
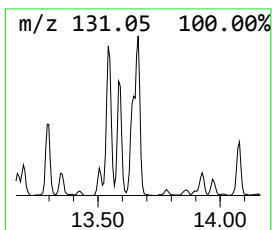
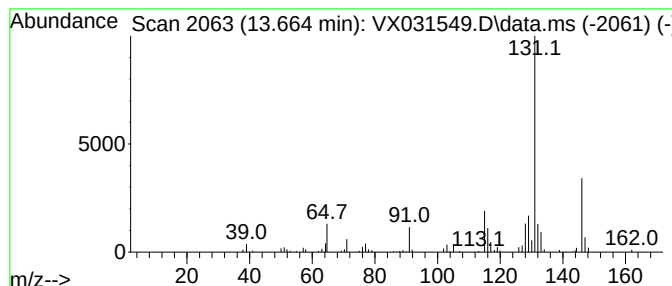
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	15.94 ug/l	106017	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91		
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	91		
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90		
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90		
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
Data File : VX031549.D
Acq On : 22 Sep 2022 04:56
Operator : JC/MD
Sample : N4642-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-3R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.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
Di-sec-Butyl ether	8.958	10.6	ug/l	72590	3	10.055	344007	50.0
Benzene, 1-ethe...	12.244	32.0	ug/l	212912	4	12.024	332499	50.0
Benzene, (1-met...	12.646	8.6	ug/l	57258	4	12.024	332499	50.0
Indan, 1-methyl-	12.701	42.4	ug/l	281852	4	12.024	332499	50.0
Benzene, (2-met...	13.134	9.3	ug/l	61743	4	12.024	332499	50.0
1H-Indene, 2,3-...	13.292	49.5	ug/l	329215	4	12.024	332499	50.0
2,2-Dimethylind...	13.542	17.8	ug/l	118228	4	12.024	332499	50.0
1H-Indene, 2,3-...	13.585	17.6	ug/l	117245	4	12.024	332499	50.0
1H-Indene, 2,3-...	13.640	16.3	ug/l	108591	4	12.024	332499	50.0
1H-Indene, 2,3-...	13.664	15.9	ug/l	106017	4	12.024	332499	50.0