

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072020\
 Data File : VX017434.D
 Acq On : 20 Jul 2020 12:40
 Operator : JC/SP
 Sample : L3329-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PW-2R

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.416	213	218	225	rVB	26311	42769	2.99%	0.288%
2	2.575	235	244	253	rBV	47834	93174	6.51%	0.627%
3	3.386	369	377	383	rBV	10002	21786	1.52%	0.147%
4	4.654	573	585	587	rBV3	7504	18625	1.30%	0.125%
5	5.105	651	659	668	rBV4	6843	17702	1.24%	0.119%
6	5.465	707	718	729	rBV2	158730	462178	32.30%	3.112%
7	5.629	735	745	759	rVB2	250515	702757	49.11%	4.732%
8	6.032	801	811	823	rBV	164440	429583	30.02%	2.893%
9	6.830	931	942	953	rBV	397175	948510	66.28%	6.387%
10	7.440	1031	1042	1049	rBV7	10108	28049	1.96%	0.189%
11	8.696	1241	1248	1256	rVB	874211	1309160	91.48%	8.815%
12	9.019	1296	1301	1308	rVB4	9596	15441	1.08%	0.104%
13	10.098	1470	1478	1486	rBV	925685	1278258	89.32%	8.607%
14	10.342	1516	1518	1524	rVB5	9790	15095	1.05%	0.102%
15	10.640	1558	1567	1571	rBV10	7132	18049	1.26%	0.122%
16	10.683	1571	1574	1582	rVB3	7428	14907	1.04%	0.100%
17	11.000	1622	1626	1630	rBV2	10350	15156	1.06%	0.102%
18	11.122	1640	1646	1651	rBV	870600	1056110	73.80%	7.111%
19	11.652	1728	1733	1743	rVB2	19908	32913	2.30%	0.222%
20	11.750	1743	1749	1753	rBV	13731	23443	1.64%	0.158%
21	11.902	1768	1774	1776	rBV3	13313	20319	1.42%	0.137%
22	11.933	1776	1779	1782	rVB	11766	15083	1.05%	0.102%
23	12.061	1794	1800	1809	rBV	1190115	1431031	100.00%	9.636%
24	12.213	1821	1825	1829	rBV	18970	23724	1.66%	0.160%
25	12.280	1829	1836	1843	rVV	674902	838880	58.62%	5.648%
26	12.372	1844	1851	1856	rVB5	21146	49689	3.47%	0.335%
27	12.445	1856	1863	1868	rBV	45275	61436	4.29%	0.414%
28	12.652	1889	1897	1900	rBV	161992	201802	14.10%	1.359%
29	12.683	1900	1902	1907	rVV	87247	105058	7.34%	0.707%
30	12.738	1907	1911	1918	rVB	326061	420122	29.36%	2.829%
31	12.799	1918	1921	1926	rVB4	15419	19383	1.35%	0.131%
32	12.878	1926	1934	1939	rVB2	44203	71341	4.99%	0.480%
33	12.963	1939	1948	1952	rBV	196192	276300	19.31%	1.860%
34	13.006	1952	1955	1961	rVB5	12512	21249	1.48%	0.143%

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Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

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 Title : SW846 8260

35	13.073	1961	1966	1970	rVB3	36861	57929	4.05%	0.390%
36	13.134	1970	1976	1979	rBV	26982	38752	2.71%	0.261%
37	13.177	1979	1983	1985	rBV2	53454	65756	4.60%	0.443%
38	13.207	1985	1988	1991	rBV	62956	60021	4.19%	0.404%
39	13.292	1997	2002	2004	rBV2	25758	35347	2.47%	0.238%
40	13.335	2004	2009	2014	rVV	375186	510300	35.66%	3.436%
41	13.390	2014	2018	2019	rVV2	27900	36270	2.53%	0.244%
42	13.408	2019	2021	2024	rVV	27888	31107	2.17%	0.209%
43	13.463	2027	2030	2035	rVB5	20501	30094	2.10%	0.203%
44	13.548	2039	2044	2046	rBV	31652	43317	3.03%	0.292%
45	13.585	2046	2050	2053	rVV	99420	135597	9.48%	0.913%
46	13.622	2053	2056	2060	rVV3	133640	192959	13.48%	1.299%
47	13.683	2060	2066	2067	rVV2	100981	161393	11.28%	1.087%
48	13.701	2067	2069	2076	rVB2	184489	279449	19.53%	1.882%
49	13.792	2081	2084	2085	rVV	29202	31737	2.22%	0.214%
50	13.817	2085	2088	2097	rVB	93614	150148	10.49%	1.011%
51	13.896	2097	2101	2105	rVB2	36362	47943	3.35%	0.323%
52	13.963	2105	2112	2116	rBV2	100431	146088	10.21%	0.984%
53	14.012	2116	2120	2124	rBV	58039	74178	5.18%	0.499%
54	14.115	2132	2137	2140	rBV	152201	185895	12.99%	1.252%
55	14.256	2155	2160	2169	rBV	176786	278854	19.49%	1.878%
56	14.359	2173	2177	2181	rBV	84106	97768	6.83%	0.658%
57	14.414	2181	2186	2191	rVB2	182869	241993	16.91%	1.629%
58	14.457	2191	2193	2199	rVB2	27738	37817	2.64%	0.255%
59	14.524	2199	2204	2208	rBV2	97062	146131	10.21%	0.984%
60	14.567	2208	2211	2215	rVB3	14815	17984	1.26%	0.121%
61	14.652	2218	2225	2227	rBV2	49021	92393	6.46%	0.622%
62	14.713	2232	2235	2240	rVB3	43781	60210	4.21%	0.405%
63	14.768	2240	2244	2249	rVB3	26417	34788	2.43%	0.234%
64	14.829	2249	2254	2257	rBV2	54320	101374	7.08%	0.683%
65	14.945	2270	2273	2279	rVB6	13772	24189	1.69%	0.163%
66	15.091	2294	2297	2302	rBV7	13358	18935	1.32%	0.127%
67	15.231	2315	2320	2327	rVB	54902	86526	6.05%	0.583%
68	15.298	2327	2331	2335	rBV7	14348	24990	1.75%	0.168%
69	15.457	2353	2357	2362	rVB	47254	66302	4.63%	0.446%
70	15.524	2364	2368	2372	rBV	37193	63075	4.41%	0.425%
71	15.566	2372	2375	2382	rVB6	27899	47561	3.32%	0.320%

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ClientSampleId :
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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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72	15.664	2385	2391	2396	rBV	191624	296397	20.71%	1.996%
73	15.786	2406	2411	2416	rBV7	14822	31953	2.23%	0.215%
74	15.865	2419	2424	2426	rVV2	50399	88560	6.19%	0.596%
75	15.896	2426	2429	2440	rVB3	64722	126815	8.86%	0.854%
76	16.042	2448	2453	2463	rVB	43184	81453	5.69%	0.548%
77	16.139	2465	2469	2477	rBV	6599	16602	1.16%	0.112%
78	16.389	2501	2510	2517	rBV	131041	258649	18.07%	1.742%
79	16.603	2539	2545	2549	rBV4	16128	37098	2.59%	0.250%
80	16.664	2551	2555	2560	rVB5	17970	30931	2.16%	0.208%
81	16.719	2560	2564	2567	rBV2	17029	28781	2.01%	0.194%

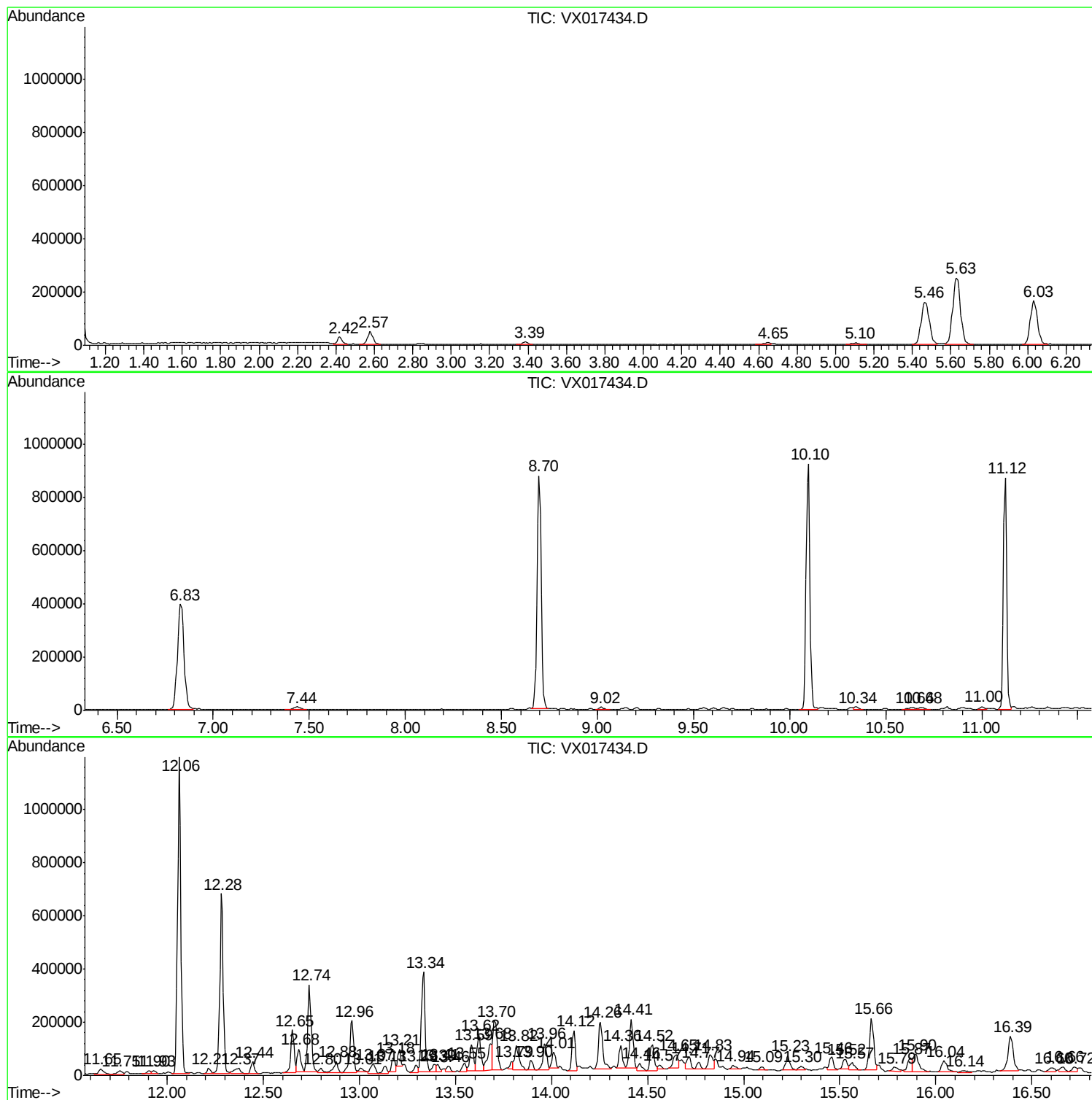
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 6 Sample Multiplier: 1

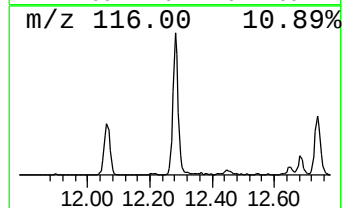
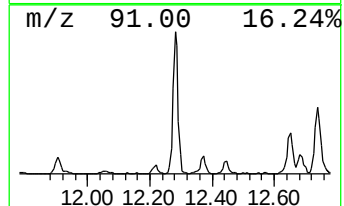
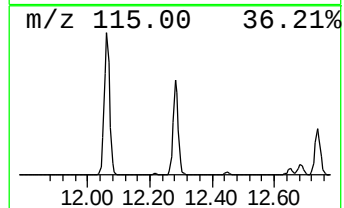
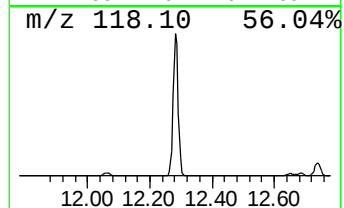
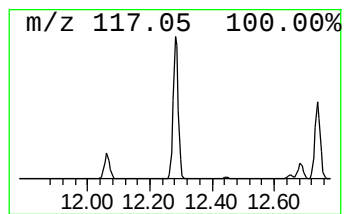
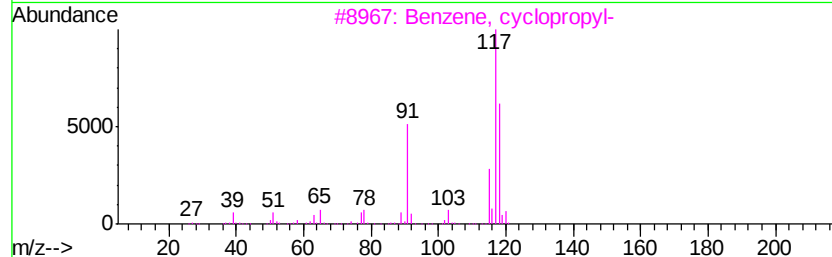
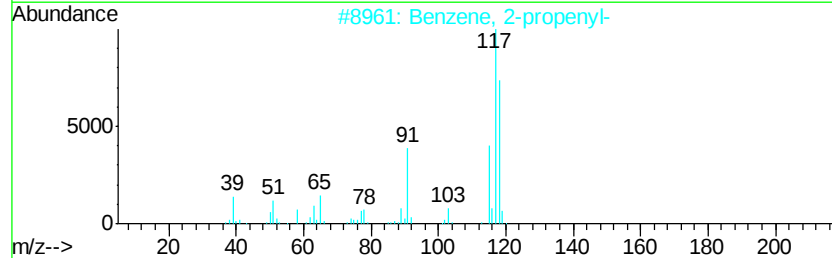
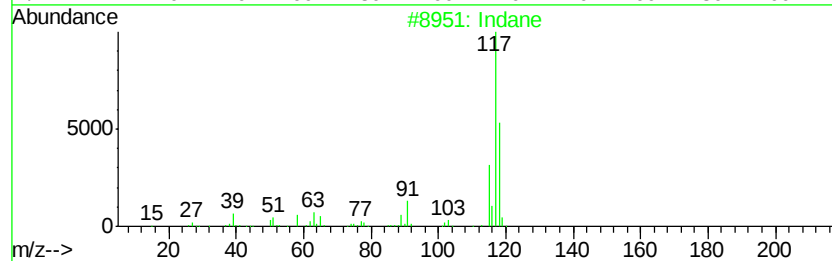
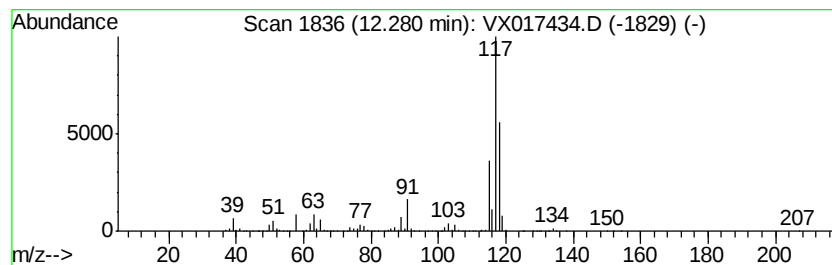
Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	29.31 ug/l	838880	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 95
2	Benzene, 2-propenyl-		118 C9H10	000300-57-2 74
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 72
4	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 72
5	7-Methylenecycloocta-1,3,5-triene		118 C9H10	002570-13-0 64



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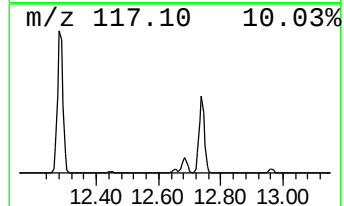
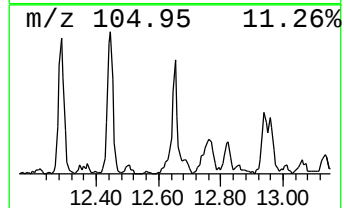
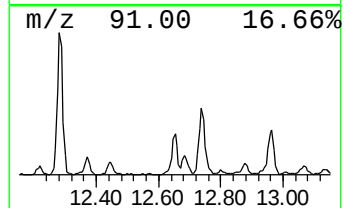
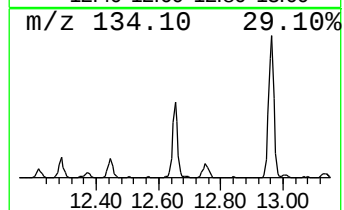
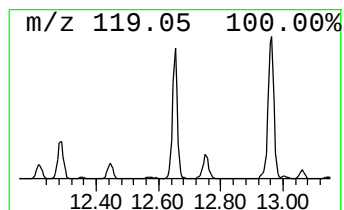
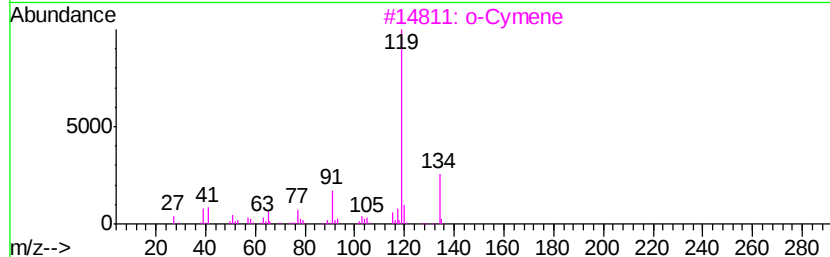
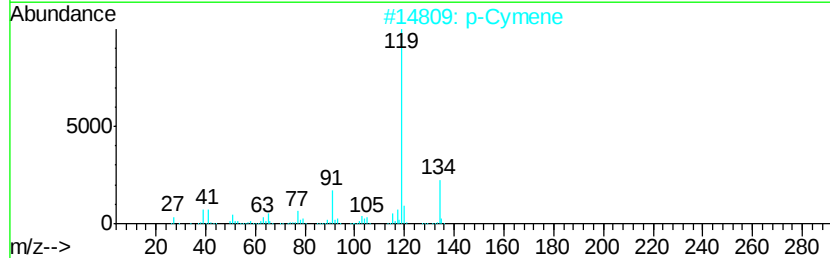
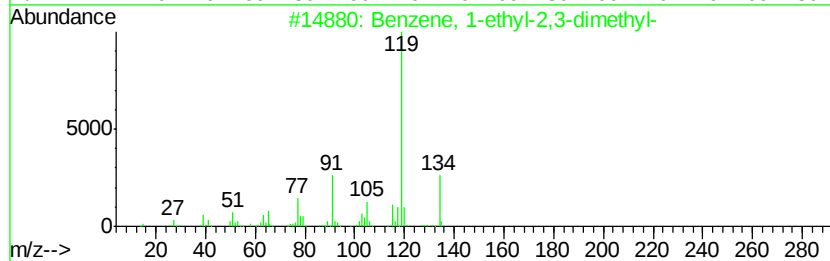
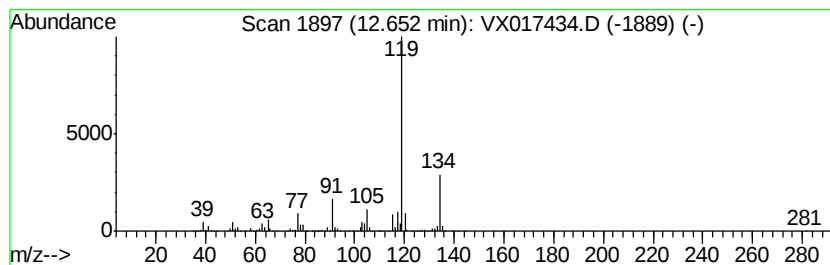
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	7.05 ug/l	201802	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		p-Cymene	134	C10H14	000099-87-6	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



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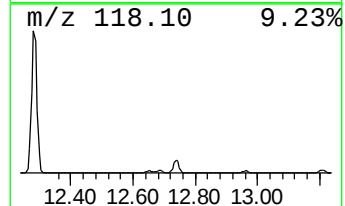
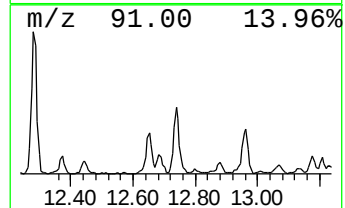
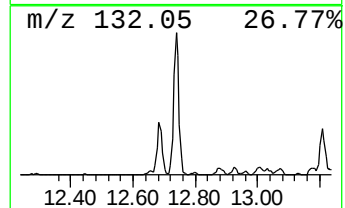
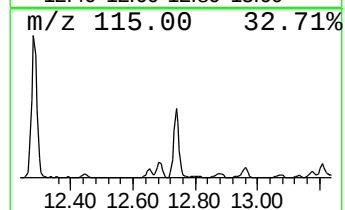
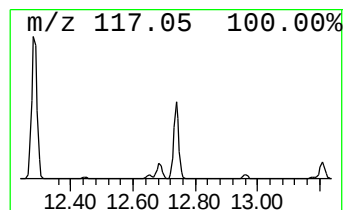
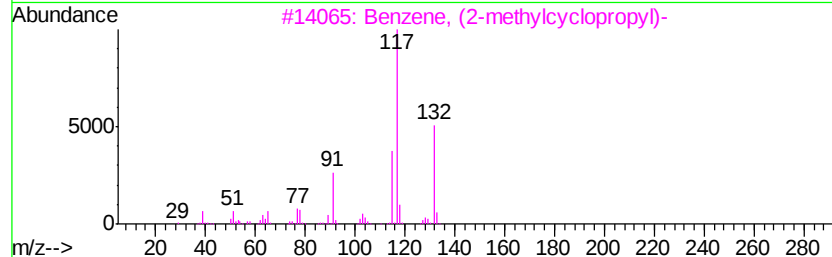
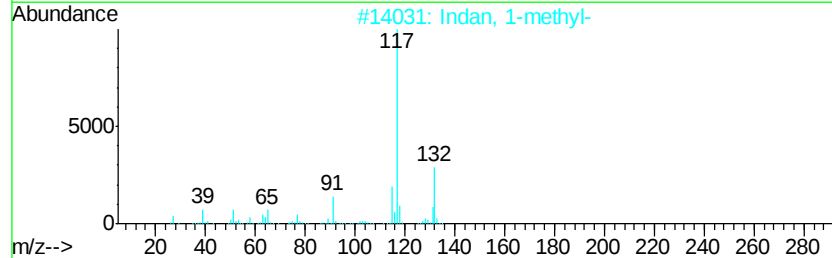
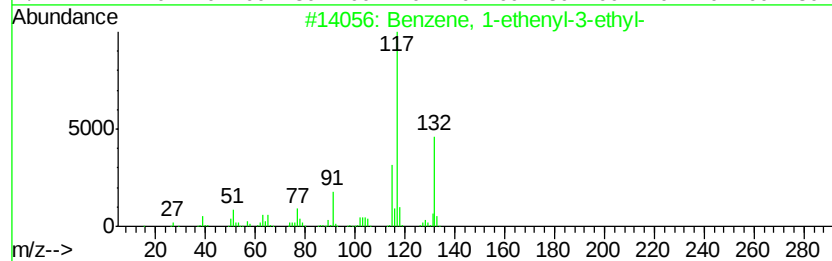
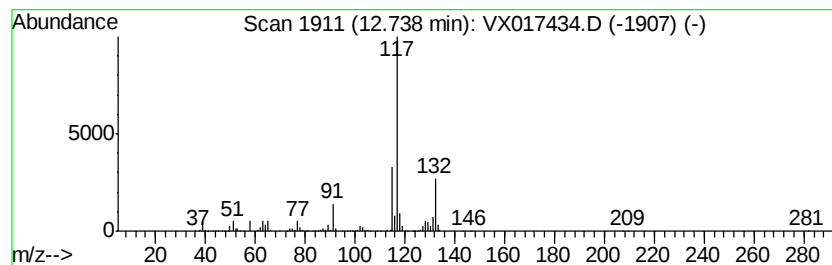
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	14.68 ug/l	420122	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	90
4		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	80
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	80



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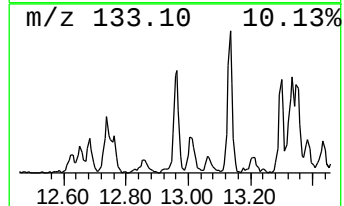
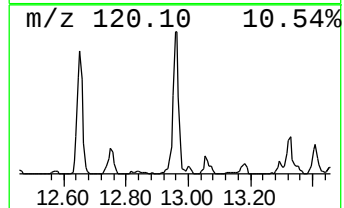
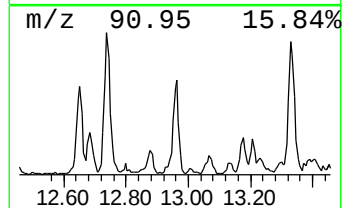
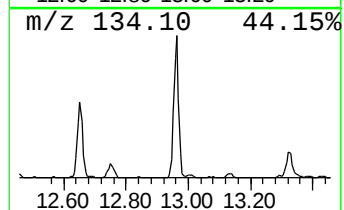
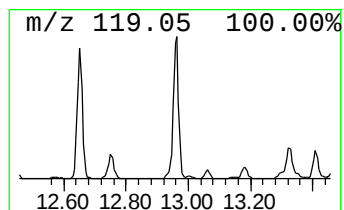
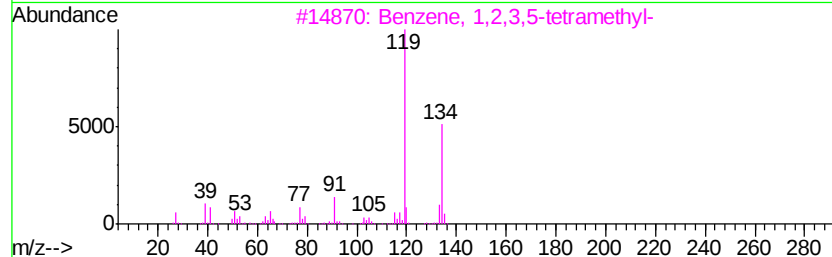
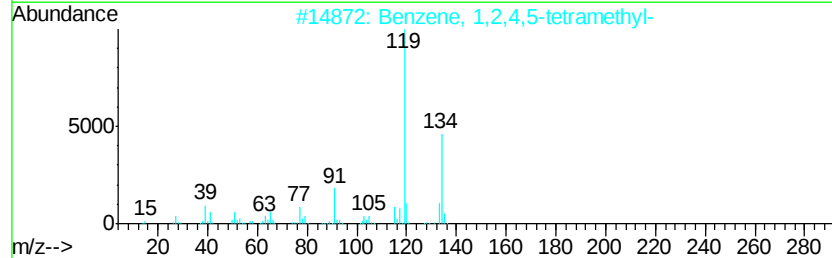
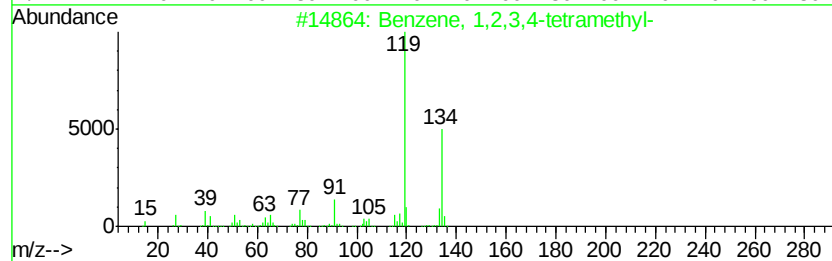
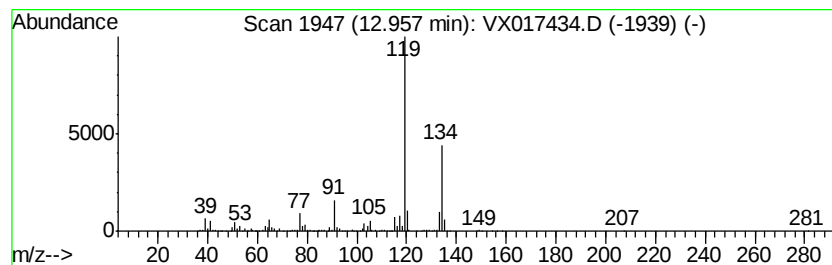
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	9.65 ug/l	276300	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97	
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96	
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94	
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91	
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PW-2R

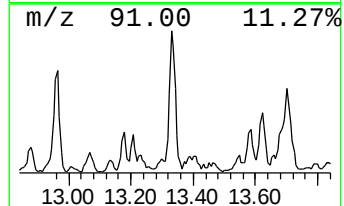
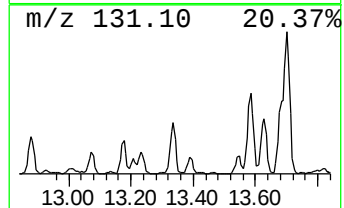
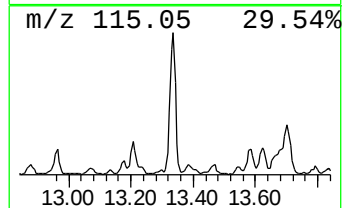
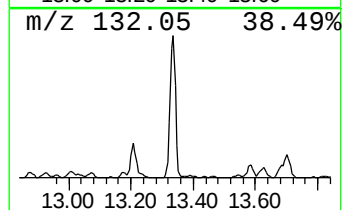
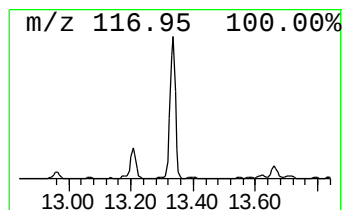
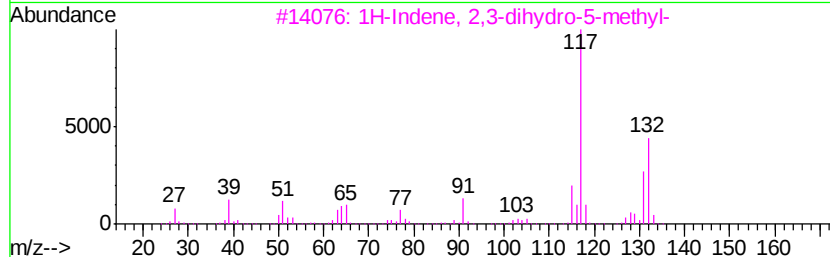
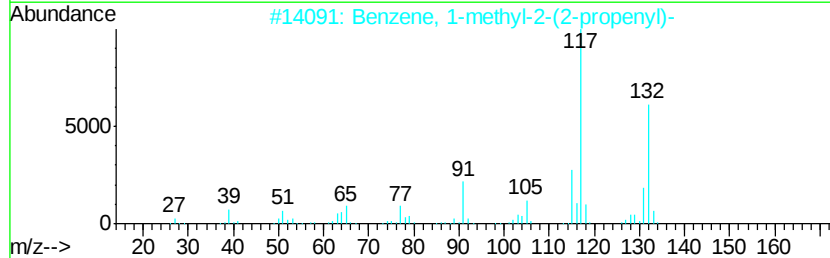
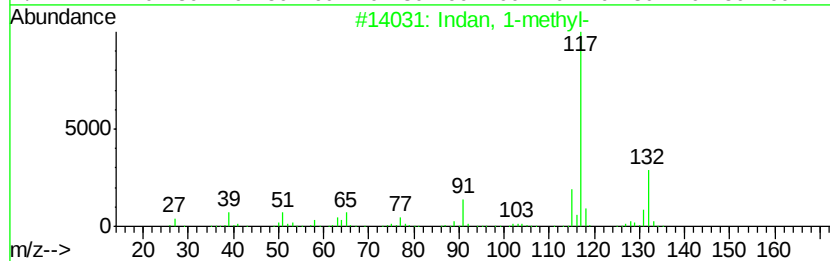
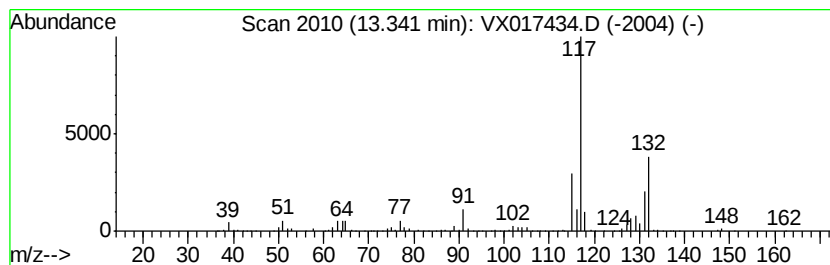
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	17.83 ug/l	510300	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	90	
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86	
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	83	
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83	
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 6 Sample Multiplier: 1

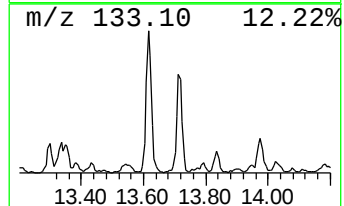
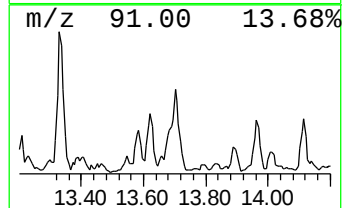
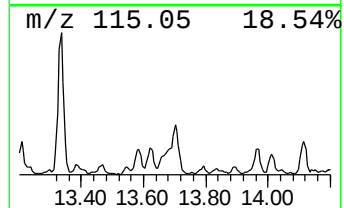
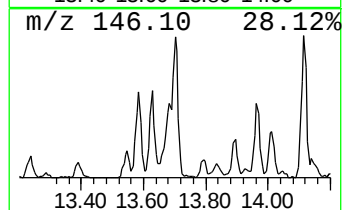
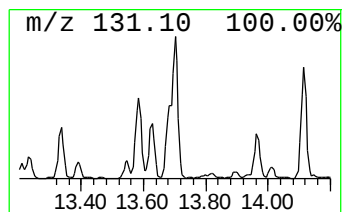
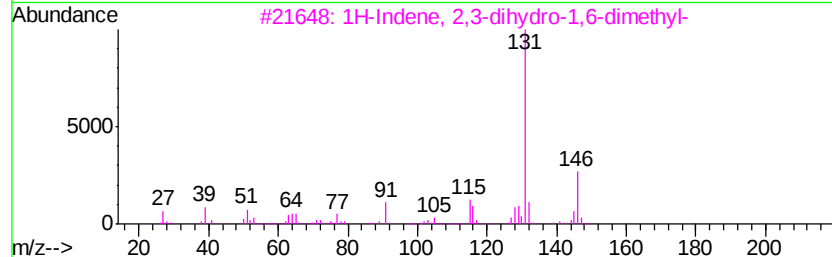
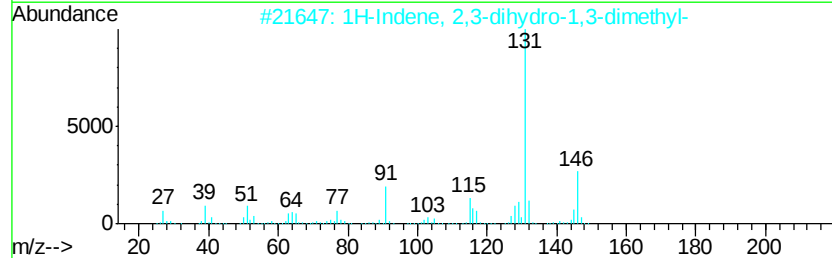
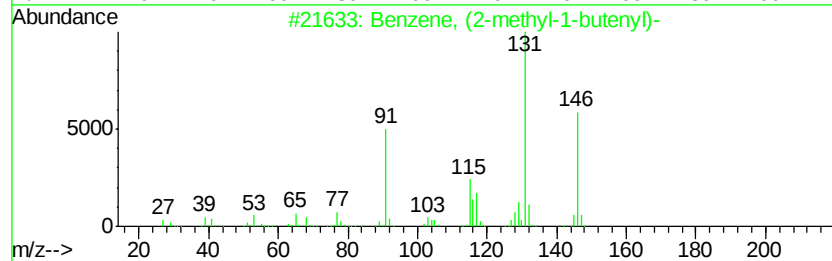
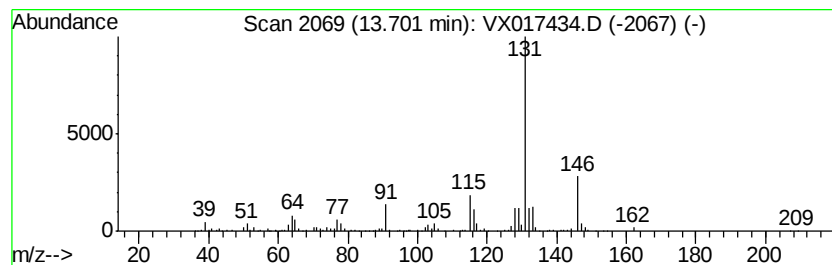
Instrument :
MSVOA_X
ClientSampleId :
PW-2R

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	9.76 ug/l	279449	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 93
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 90
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 90
4		Naphthalene, 1,2,3,4-tetrahydro...	146 C11H14	001559-81-5 90
5		1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PW-2R

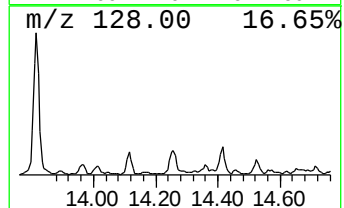
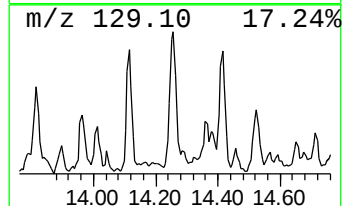
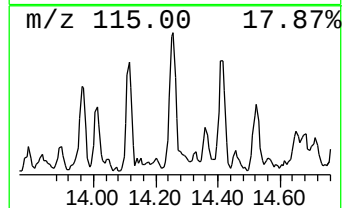
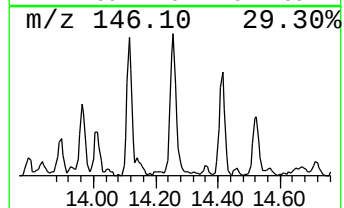
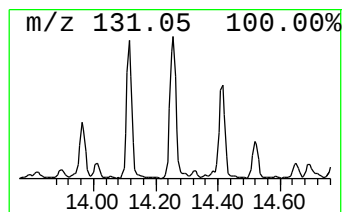
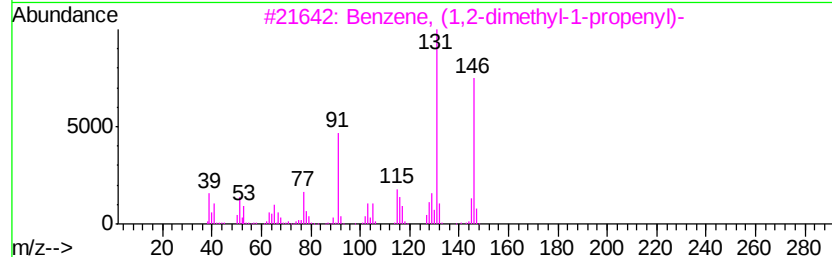
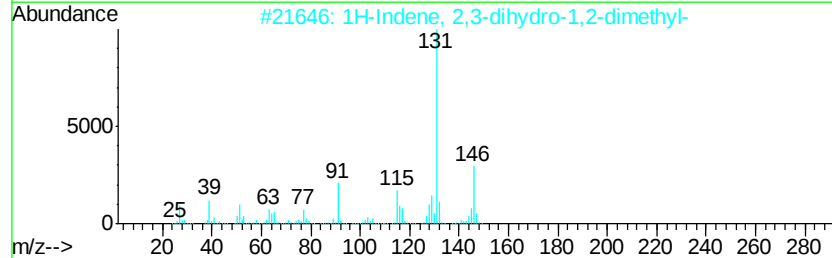
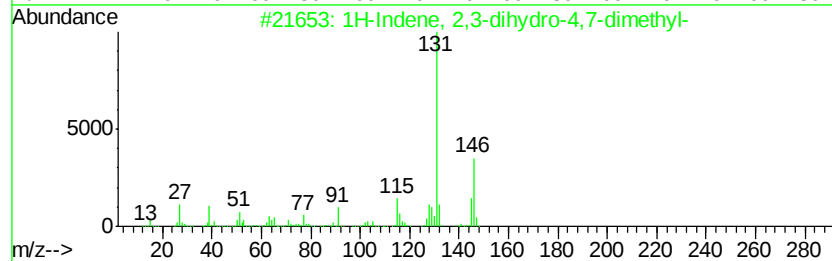
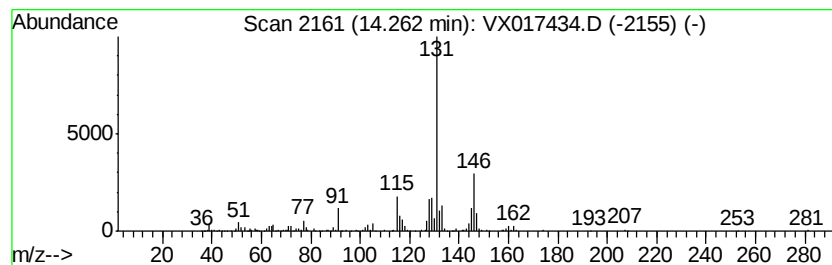
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	9.74 ug/l	278854	1,4-Dichlorobenzene-d4	12.06

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-1,2-dimet...	146	C11H14		017057-82-8	93
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14		000769-57-3	87
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14		053172-84-2	83
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14		004489-84-3	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PW-2R

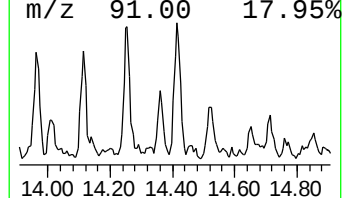
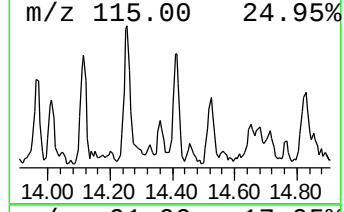
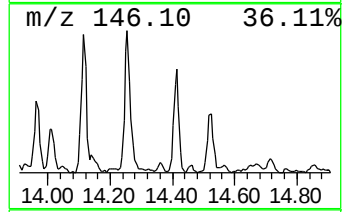
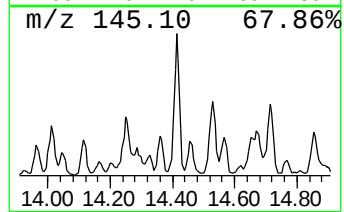
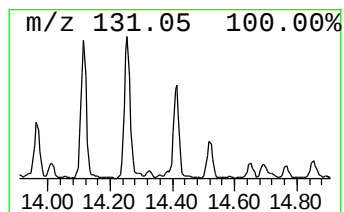
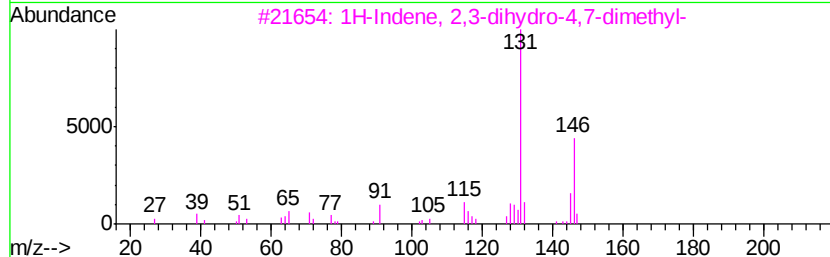
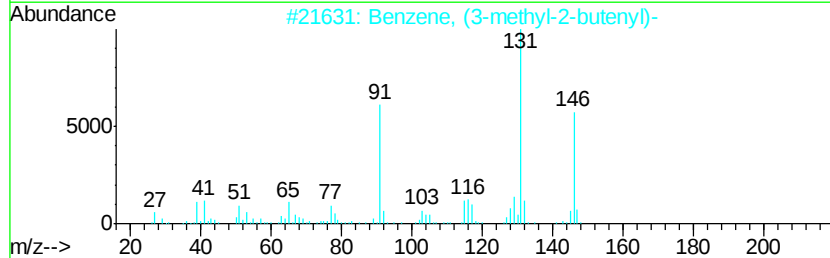
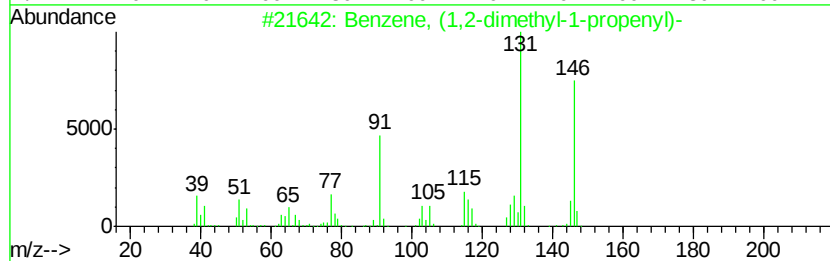
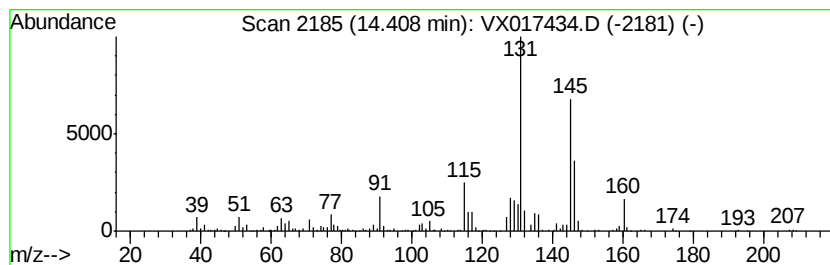
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	8.46 ug/l	241993	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	50
2		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	50
3		1H-Indene, 2,3-dihydro-4,7-dimethyl-	146	C11H14	006682-71-9	49
4		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	45
5		Benzene, 1-methyl-3-(1-methyl-2-propenyl)-	146	C11H14	052161-57-6	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PW-2R

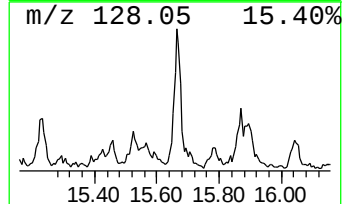
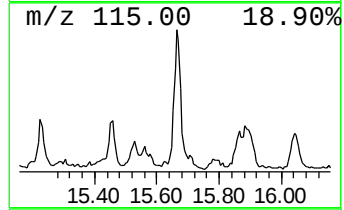
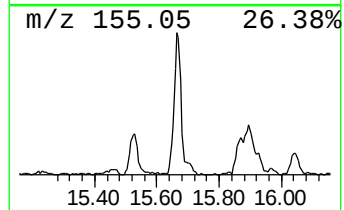
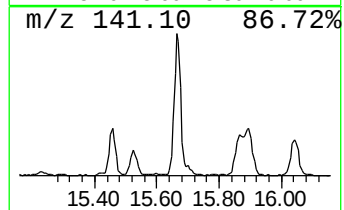
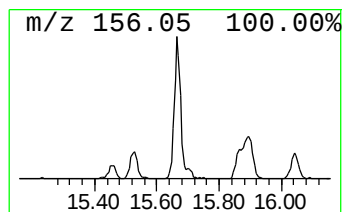
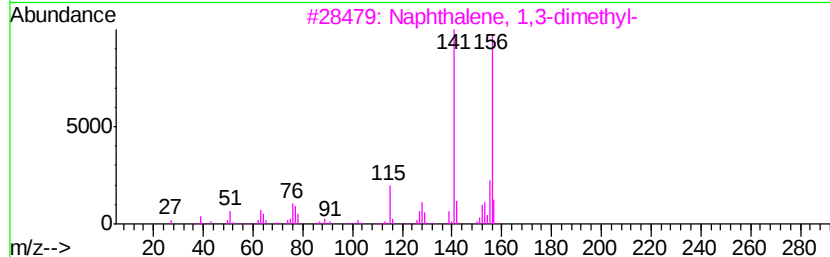
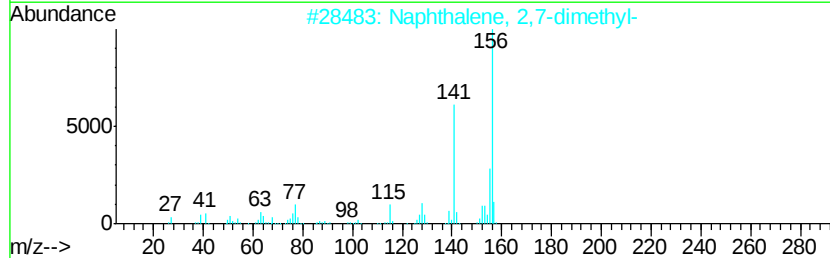
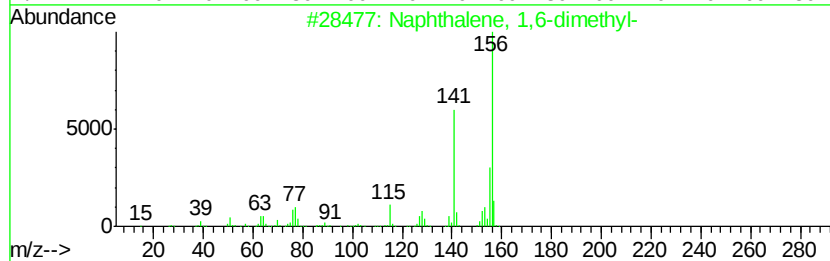
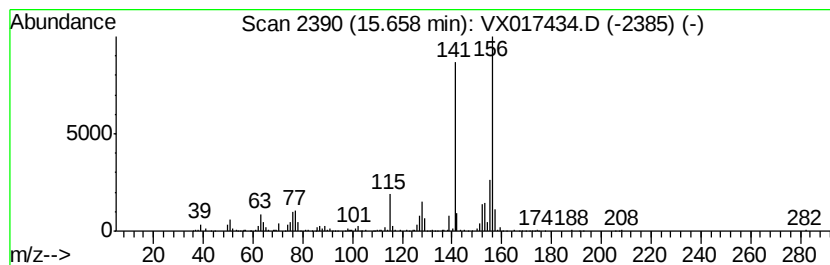
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	10.36 ug/l	296397	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072020\
Data File : VX017434.D
Acq On : 20 Jul 2020 12:40
Operator : JC/SP
Sample : L3329-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PW-2R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071320W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.28	29.3	ug/l	838880	4	12.06	1431030	50.0
Benzene, 1-ethyl-...	12.65	7.0	ug/l	201802	4	12.06	1431030	50.0
Benzene, 1-etheny...	12.74	14.7	ug/l	420122	4	12.06	1431030	50.0
Benzene, 1,2,3,4-...	12.96	9.7	ug/l	276300	4	12.06	1431030	50.0
Indan, 1-methyl-	13.34	17.8	ug/l	510300	4	12.06	1431030	50.0
Benzene, (2-methy...	13.70	9.8	ug/l	279449	4	12.06	1431030	50.0
1H-Indene, 2,3-di...	14.26	9.7	ug/l	278854	4	12.06	1431030	50.0
Benzene, (1,2-dim...	14.41	8.5	ug/l	241993	4	12.06	1431030	50.0
Naphthalene, 1,6-...	15.66	10.4	ug/l	296397	4	12.06	1431030	50.0