

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW082219\
 Data File : VW012109.D
 Acq On : 22 Aug 2019 14:34
 Operator : SY/VA
 Sample : K4086-01
 Misc : 4.01G/5ML/MSVOA W/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 NB-08-082019

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_W\METHOD\82W072719S.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	2.148	36	40	46	rBV	13499	24825	1.98%	0.305%
2	4.915	482	494	511	rBV2	41446	145483	11.60%	1.788%
3	5.940	651	662	673	rBV3	11003	31764	2.53%	0.390%
4	7.537	913	924	935	rBV3	6694	19940	1.59%	0.245%
5	7.884	971	981	986	rBV	163995	388537	30.99%	4.775%
6	7.945	986	991	1003	rVB	304741	671611	53.56%	8.254%
7	8.305	1040	1050	1060	rBV2	181947	402152	32.07%	4.943%
8	8.841	1127	1138	1151	rBV	483928	947072	75.53%	11.640%
9	10.323	1373	1381	1389	rBV	706096	1253837	100.00%	15.410%
10	10.384	1389	1391	1397	rVB	8190	12653	1.01%	0.156%
11	10.707	1439	1444	1449	rVB2	10479	18084	1.44%	0.222%
12	11.634	1588	1596	1604	rBV	546844	940976	75.05%	11.565%
13	12.164	1672	1683	1693	rBV7	5722	15958	1.27%	0.196%
14	12.621	1750	1758	1766	rBV	378668	644811	51.43%	7.925%
15	12.938	1805	1810	1817	rVB4	10116	20509	1.64%	0.252%
16	13.249	1857	1861	1867	rVB	9529	14872	1.19%	0.183%
17	13.560	1905	1912	1922	rBV	381237	625417	49.88%	7.686%
18	13.755	1940	1944	1947	rBV3	11932	18979	1.51%	0.233%
19	13.792	1947	1950	1959	rVB3	14416	30274	2.41%	0.372%
20	13.877	1959	1964	1971	rBV3	15134	27061	2.16%	0.333%
21	14.011	1982	1986	1988	rBV2	9112	13121	1.05%	0.161%
22	14.072	1993	1996	1998	rBV2	9977	13781	1.10%	0.169%
23	14.206	2013	2018	2023	rVB2	21302	33691	2.69%	0.414%
24	14.334	2034	2039	2044	rBV5	9965	18072	1.44%	0.222%
25	14.420	2050	2053	2056	rVV	14064	20668	1.65%	0.254%
26	14.456	2056	2059	2063	rVV	19623	26187	2.09%	0.322%
27	14.505	2063	2067	2069	rVV2	12406	17045	1.36%	0.209%
28	14.542	2070	2073	2081	rVB3	12021	20434	1.63%	0.251%
29	14.639	2081	2089	2093	rBV3	15684	34874	2.78%	0.429%
30	14.694	2093	2098	2100	rBV2	27592	46808	3.73%	0.575%
31	14.725	2100	2103	2109	rVB3	33410	57292	4.57%	0.704%
32	14.810	2109	2117	2122	rBV6	67357	158235	12.62%	1.945%
33	14.859	2122	2125	2133	rVB3	12220	23078	1.84%	0.284%
34	14.962	2136	2142	2148	rBV	46995	78973	6.30%	0.971%

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35	15.066	2149	2159	2163	rBV3	31048	77346	6.17%	0.951%
36	15.109	2163	2166	2170	rVV	16045	22879	1.82%	0.281%
37	15.182	2171	2178	2184	rVB3	35078	80761	6.44%	0.993%
38	15.255	2186	2190	2197	rVB7	8085	16591	1.32%	0.204%
39	15.334	2197	2203	2205	rBV5	13311	26093	2.08%	0.321%
40	15.426	2212	2218	2222	rBV2	37673	72658	5.79%	0.893%
41	15.499	2222	2230	2236	rVV	76112	194843	15.54%	2.395%
42	15.554	2236	2239	2249	rVB3	27499	57165	4.56%	0.703%
43	15.657	2249	2256	2264	rBV4	38173	87297	6.96%	1.073%
44	15.828	2273	2284	2286	rBV6	28285	74441	5.94%	0.915%
45	15.865	2286	2290	2299	rVB	86348	163702	13.06%	2.012%
46	15.956	2300	2305	2310	rBV6	12151	22914	1.83%	0.282%
47	16.035	2310	2318	2323	rBV4	20790	53895	4.30%	0.662%
48	16.182	2333	2342	2348	rBV	54186	127884	10.20%	1.572%
49	16.377	2365	2374	2380	rBV3	54809	129721	10.35%	1.594%
50	16.444	2380	2385	2390	rBV3	20314	42239	3.37%	0.519%
51	16.651	2411	2419	2426	rBV8	20347	69104	5.51%	0.849%

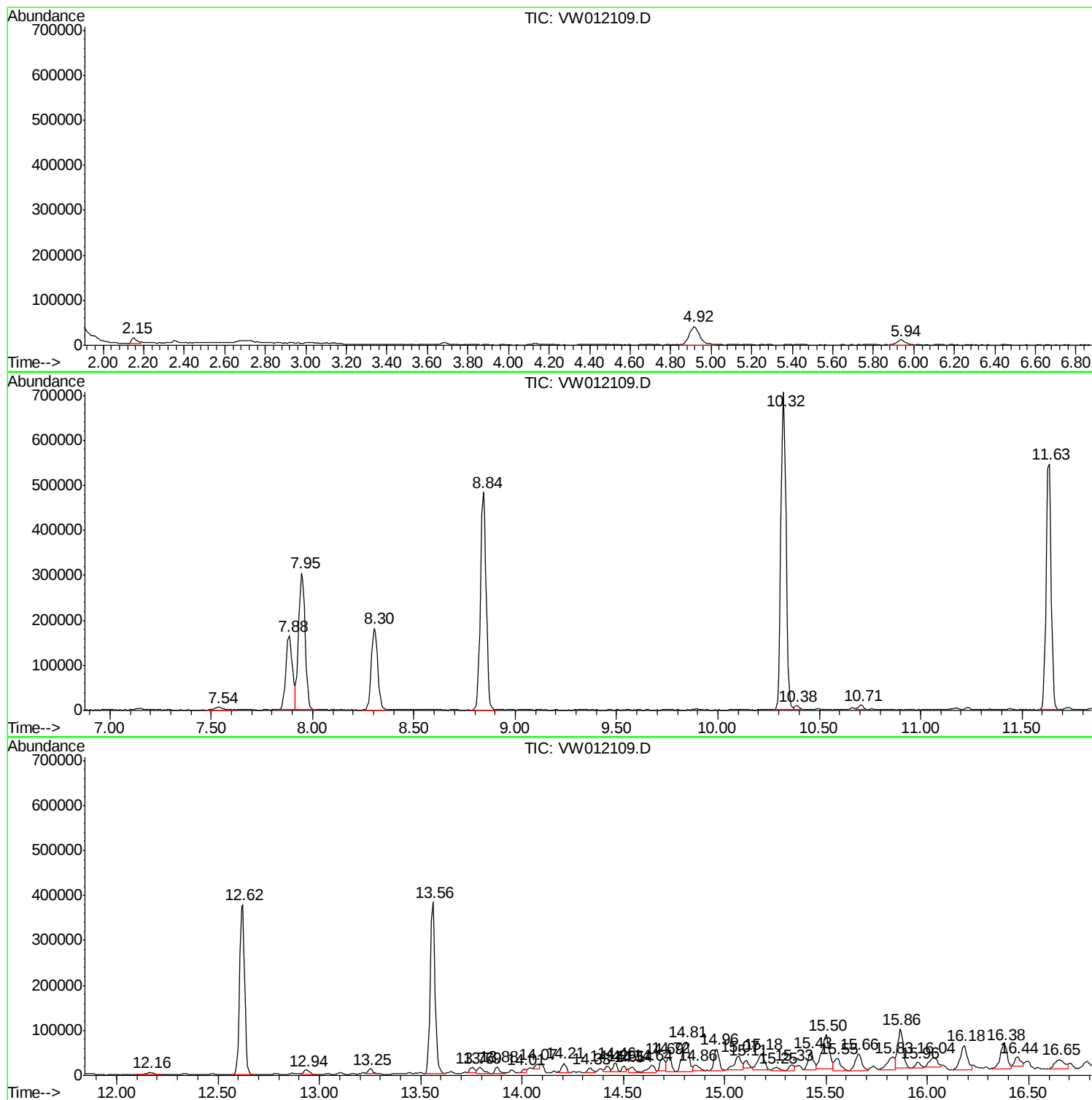
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.M
Quant Title : SW846 8260

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



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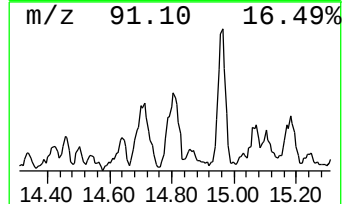
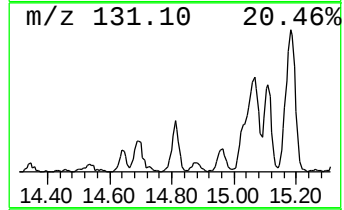
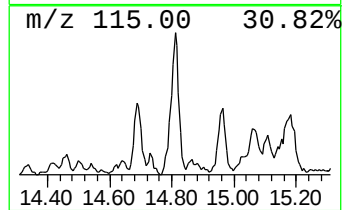
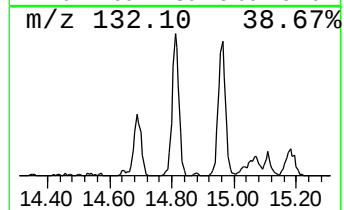
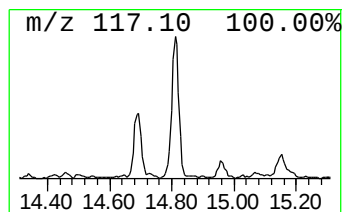
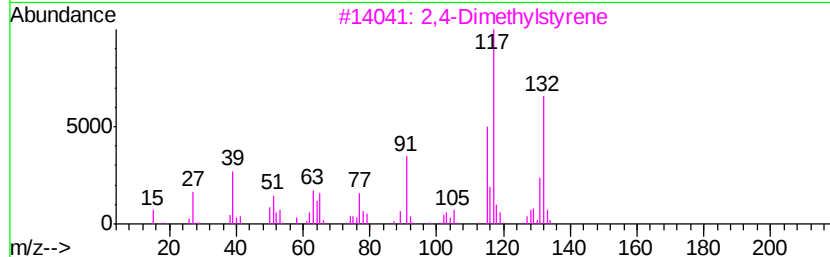
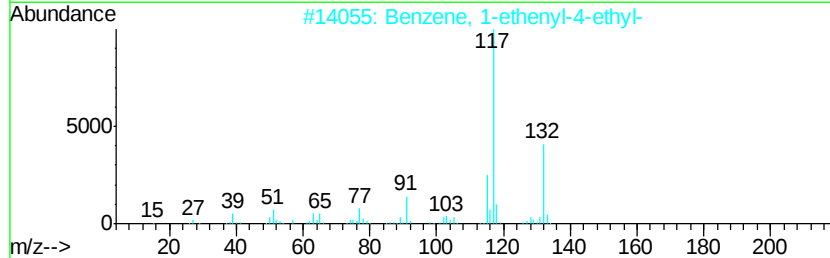
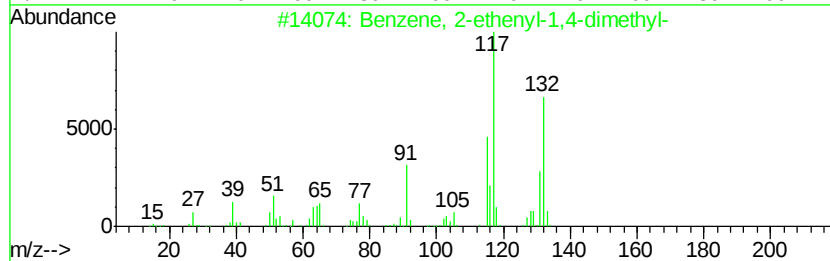
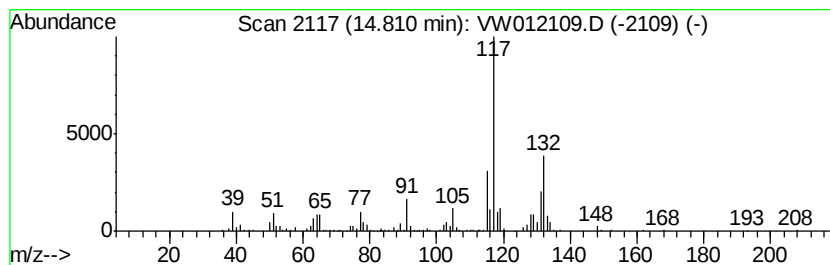
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	12.65 ug/l	158235	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	92
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90



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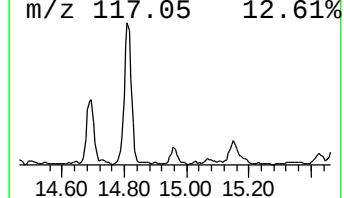
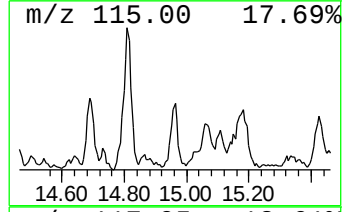
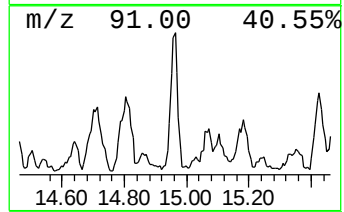
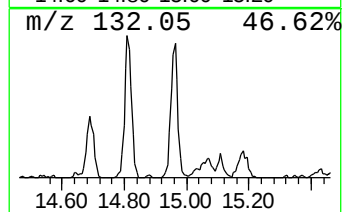
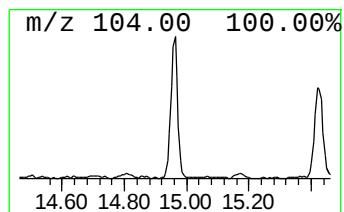
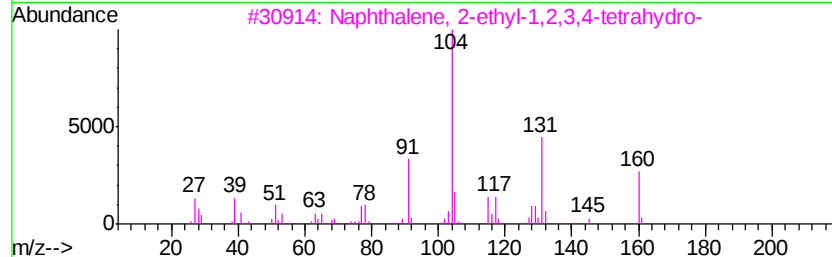
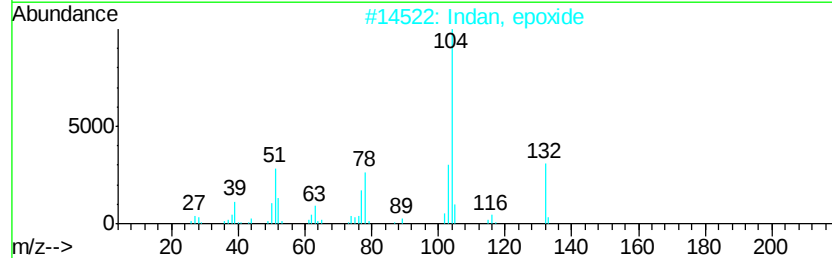
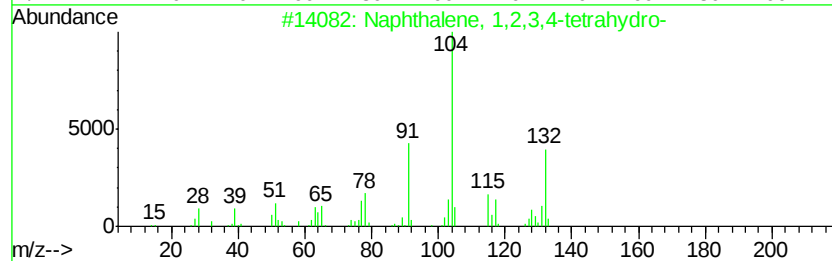
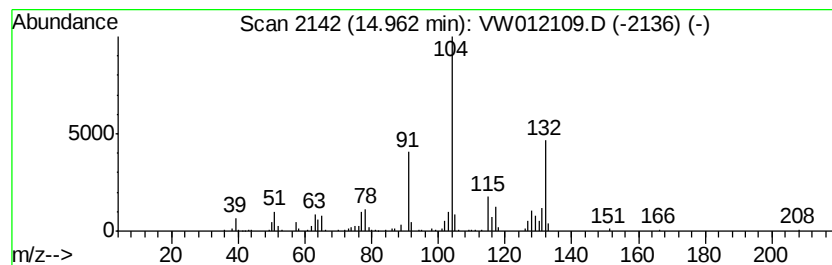
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	6.31 ug/l	78973	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		Indan, epoxide	132	C9H8O	000768-22-9	58
3		Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	53
4		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	43
5		Styrene	104	C8H8	000100-42-5	38



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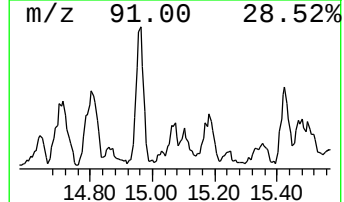
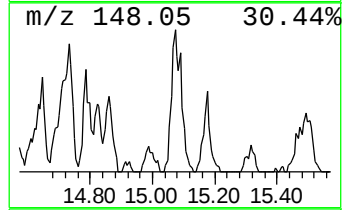
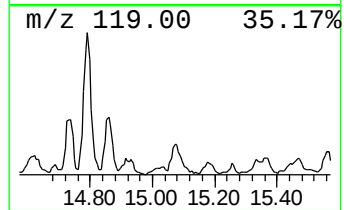
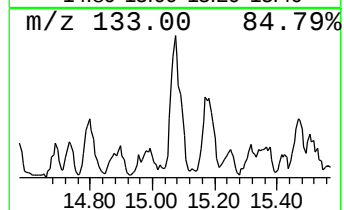
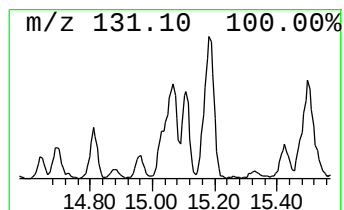
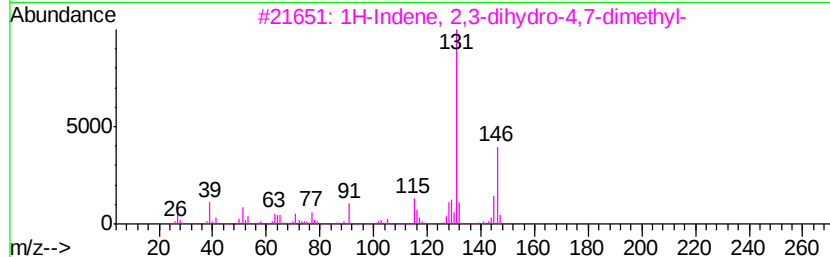
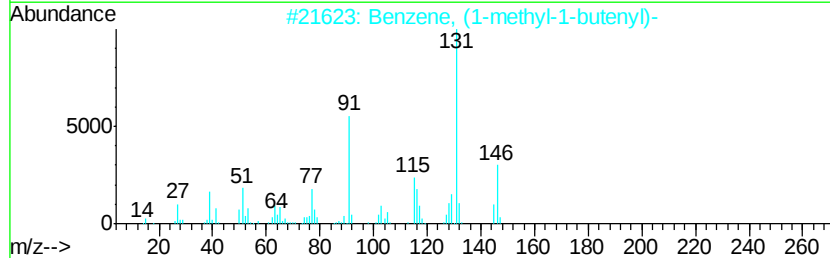
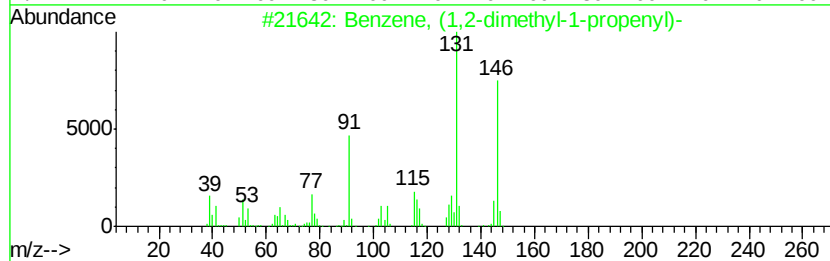
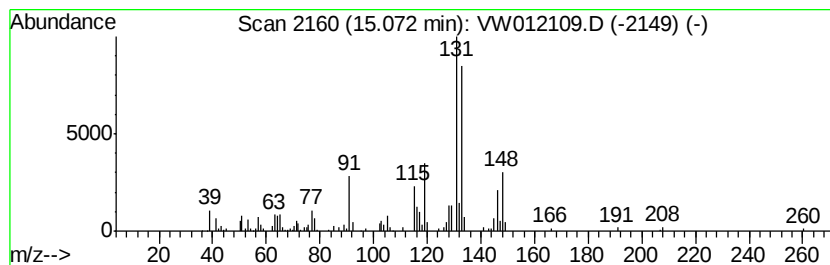
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	6.18 ug/l	77346	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	93
3		1H-Indene, 2,3-dihydro-4,7-dimethyl-	146	C11H14	006682-71-9	90
4		1H-Indene, 2,3-dihydro-1,2-dimethyl-	146	C11H14	017057-82-8	90
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83



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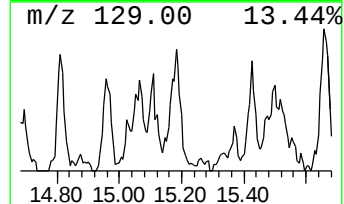
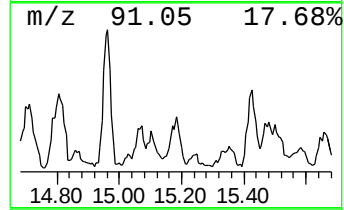
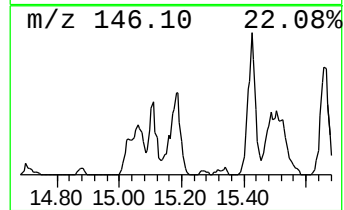
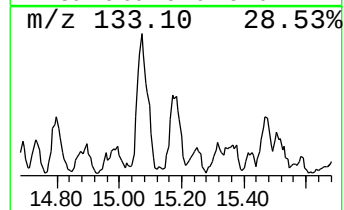
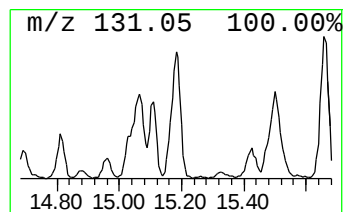
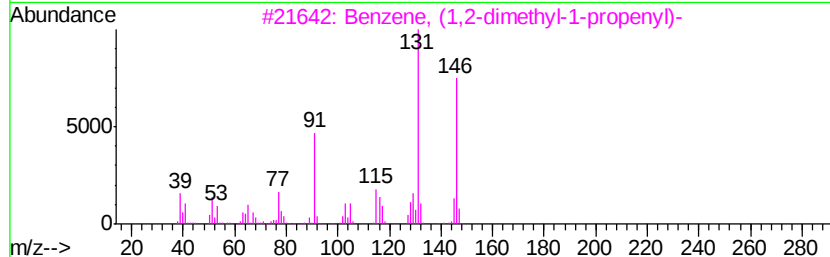
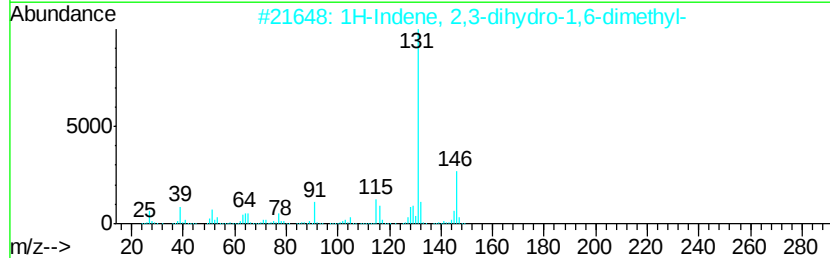
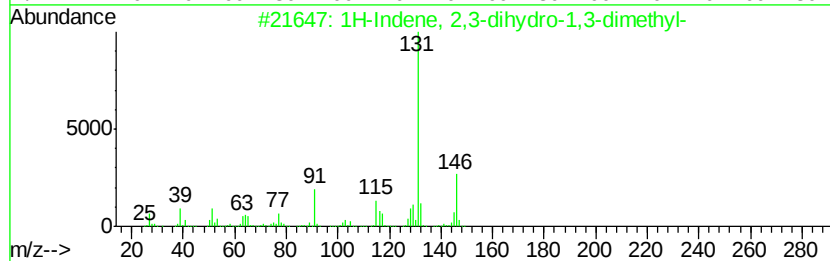
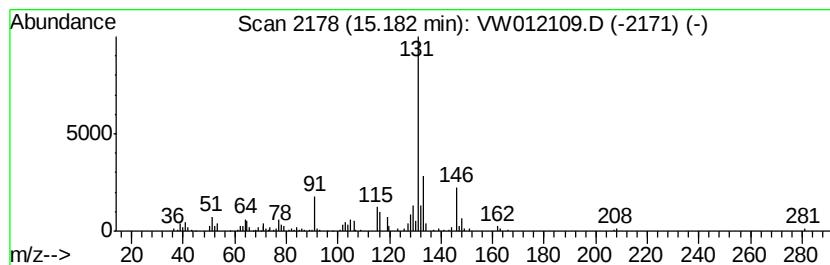
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.18	6.46 ug/l	80761	1,4-Dichlorobenzene-d4		13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14		004175-53-5	94
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14		017059-48-2	94
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14		000769-57-3	93
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14		004489-84-3	89
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14		020836-11-7	76



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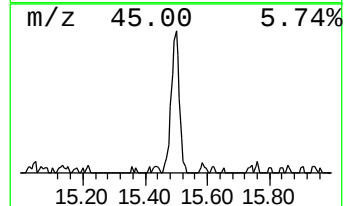
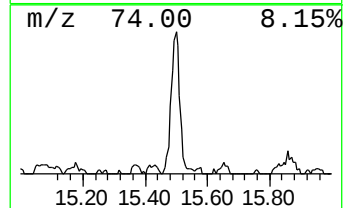
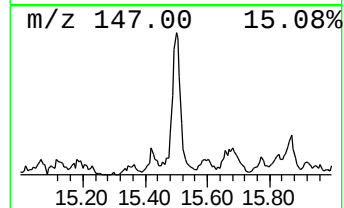
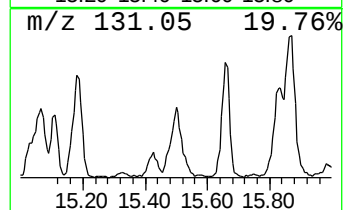
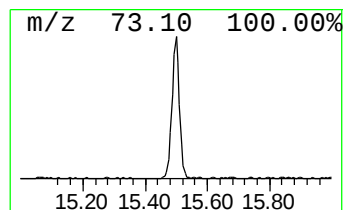
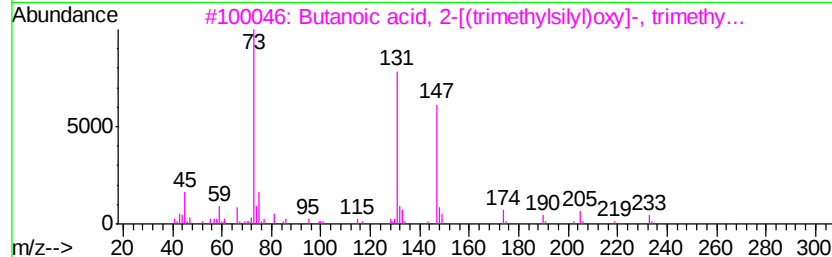
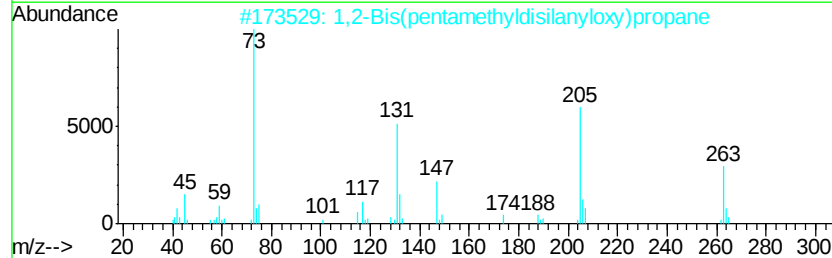
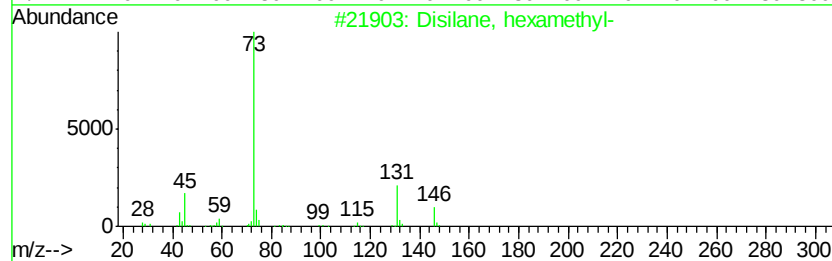
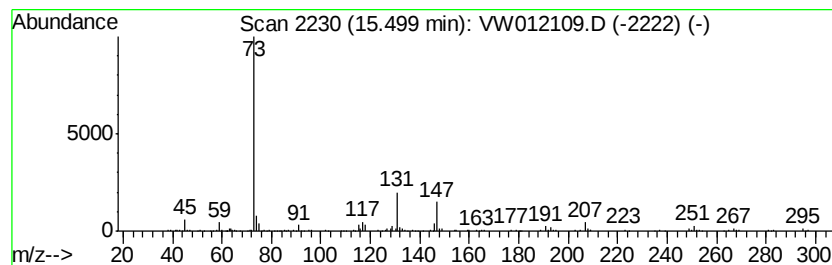
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Quant Title : SW846 8260

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

Peak Number 5 Disilane, hexamethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.50	15.58 ug/l	194843	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Disilane, hexamethyl-	146	C6H18Si2	001450-14-2	43
2		1,2-Bis(pentamethylidisilanyloxy)...	336	C13H36O2Si4	1000216-98-7	42
3		Butanoic acid, 2-[(trimethylsilyl)...	248	C10H24O3Si2	055133-93-2	40
4		2-Methyl-3,5-heptadiene, 7-trime...	358	C17H38O2Si3	1000196-08-4	38
5		Propane, 2-methyl-1,2-bis(trimet...	234	C10H26O2Si2	099875-05-5	33



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082219\
Data File : VW012109.D
Acq On : 22 Aug 2019 14:34
Operator : SY/VA
Sample : K4086-01
Misc : 4.01G/5ML/MSVOA W/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
NB-08-082019

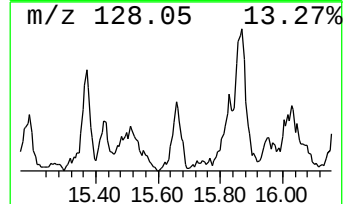
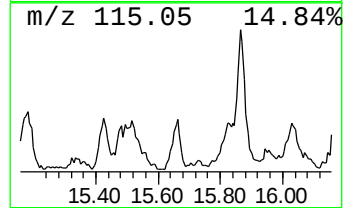
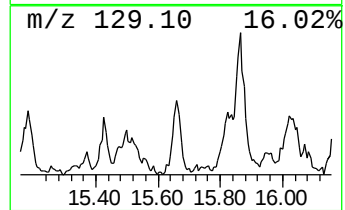
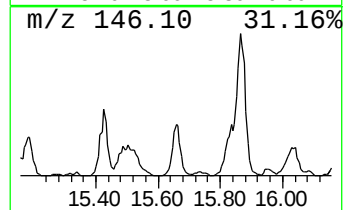
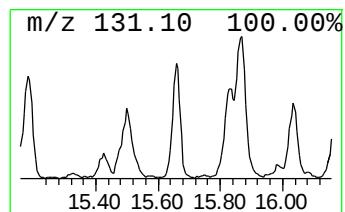
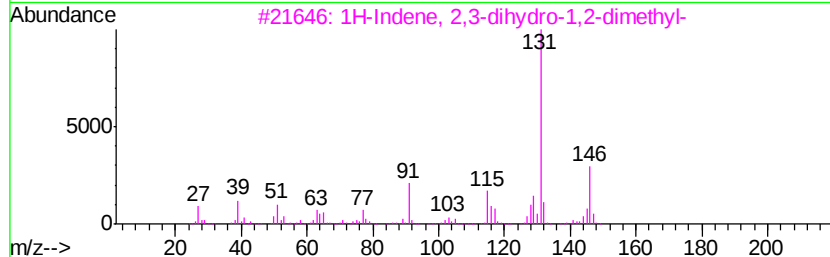
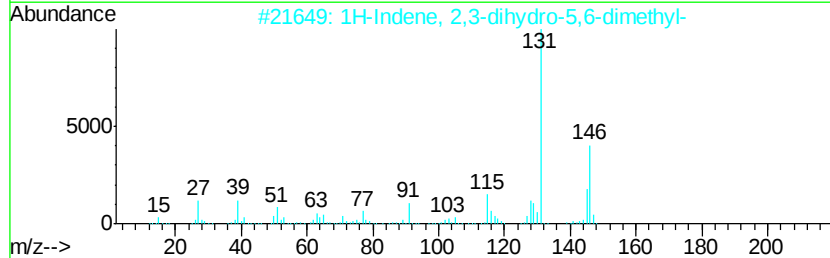
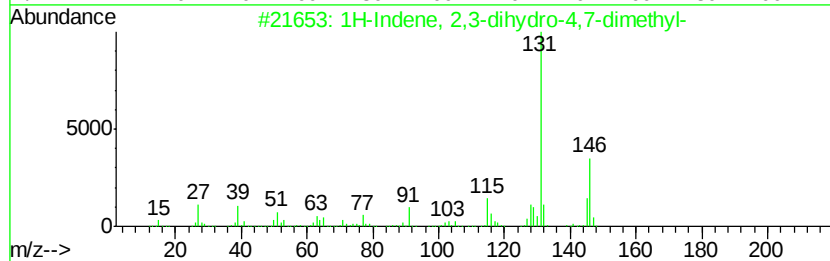
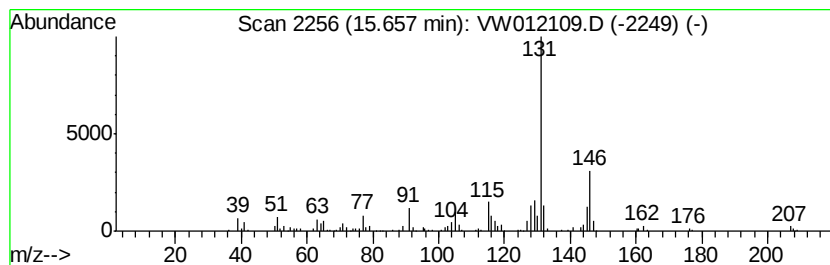
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	6.98 ug/l	87297	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	95
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	95
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082219\
Data File : VW012109.D
Acq On : 22 Aug 2019 14:34
Operator : SY/VA
Sample : K4086-01
Misc : 4.01G/5ML/MSVOA W/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
NB-08-082019

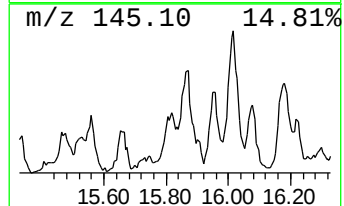
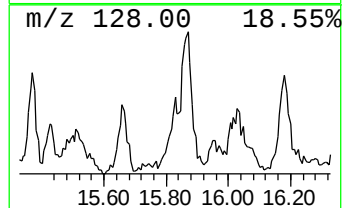
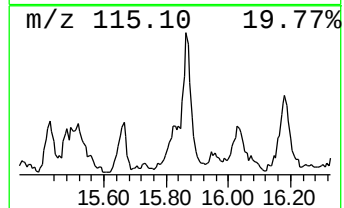
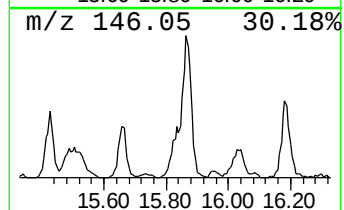
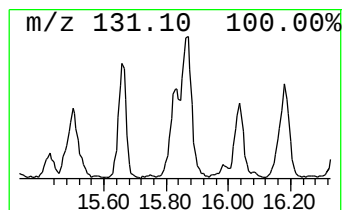
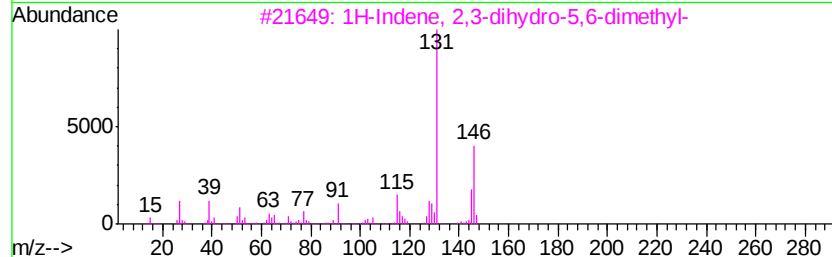
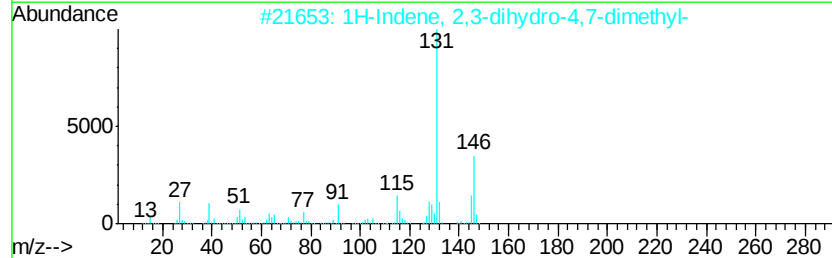
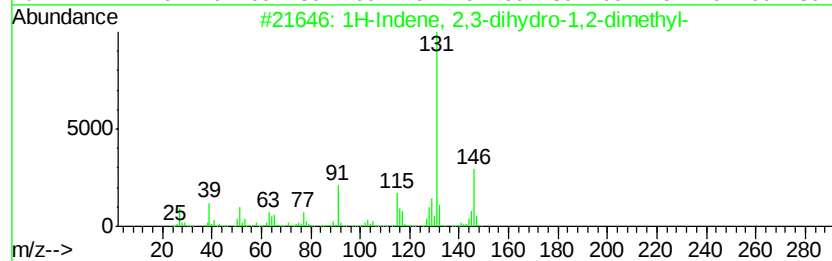
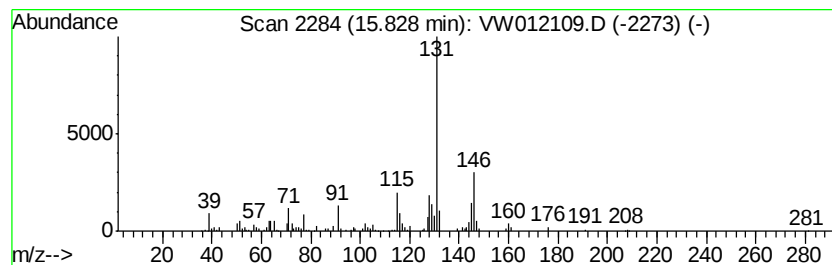
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.83	5.95 ug/l	74441	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	87
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	87
5		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082219\
Data File : VW012109.D
Acq On : 22 Aug 2019 14:34
Operator : SY/VA
Sample : K4086-01
Misc : 4.01G/5ML/MSVOA W/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
NB-08-082019

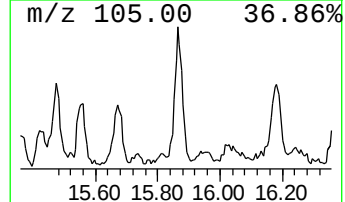
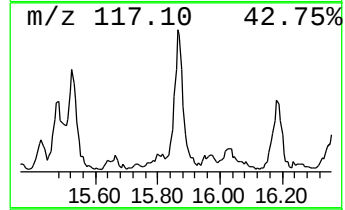
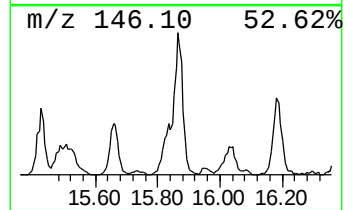
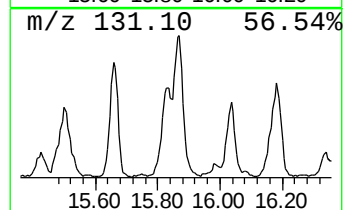
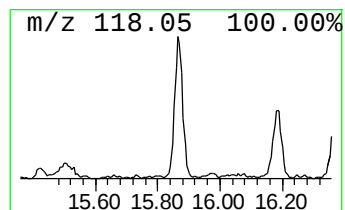
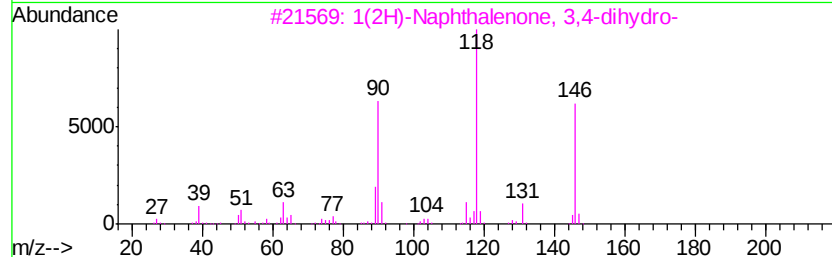
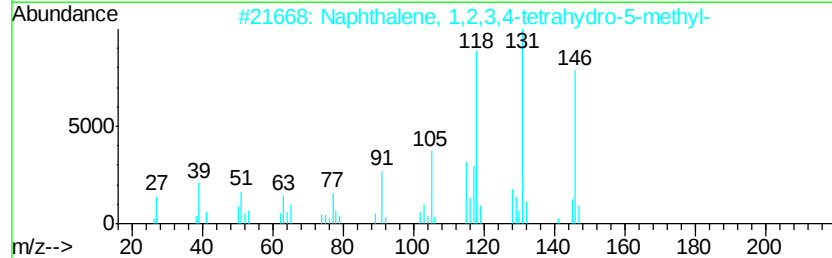
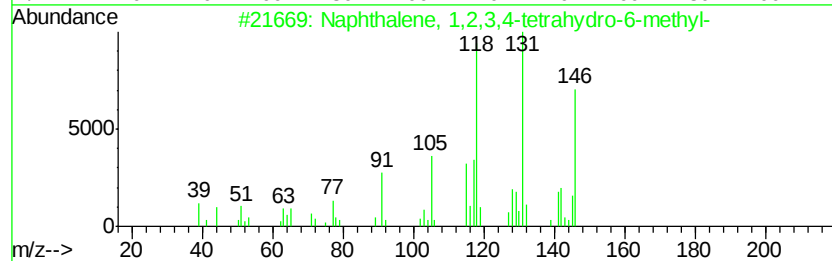
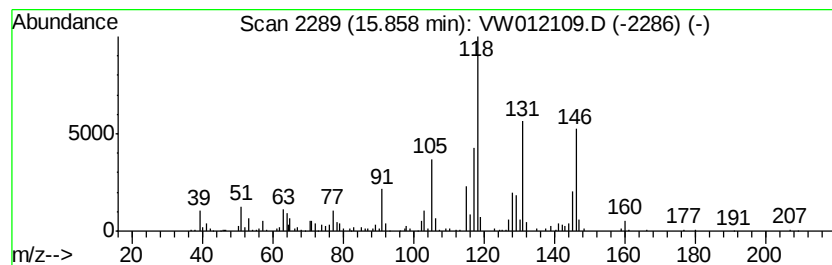
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	13.09 ug/l	163702	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	53
3		1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	43
4		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	41
5		2(1H)-Quinoxalinone	146	C8H6N2O	001196-57-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW082219\
Data File : VW012109.D
Acq On : 22 Aug 2019 14:34
Operator : SY/VA
Sample : K4086-01
Misc : 4.01G/5ML/MSVOA W/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
NB-08-082019

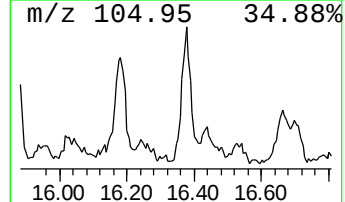
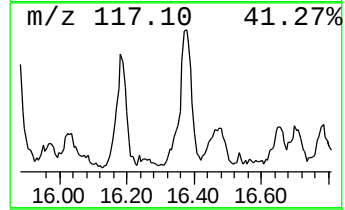
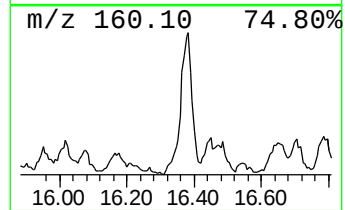
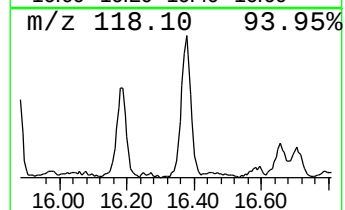
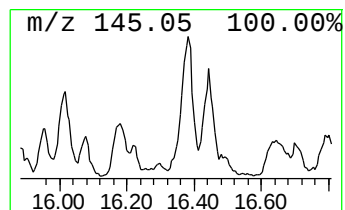
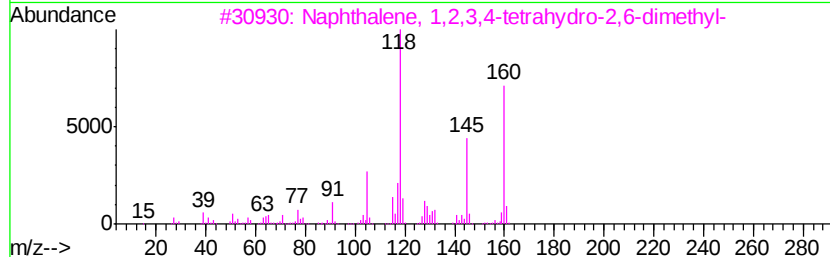
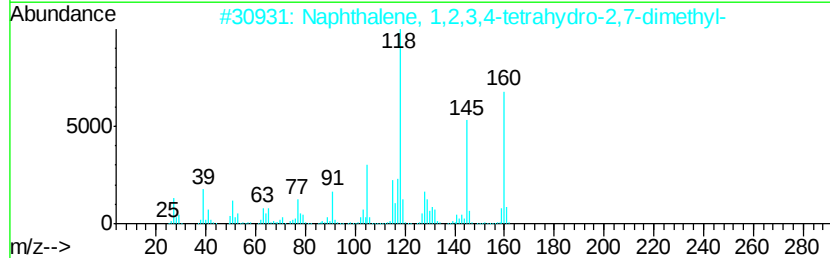
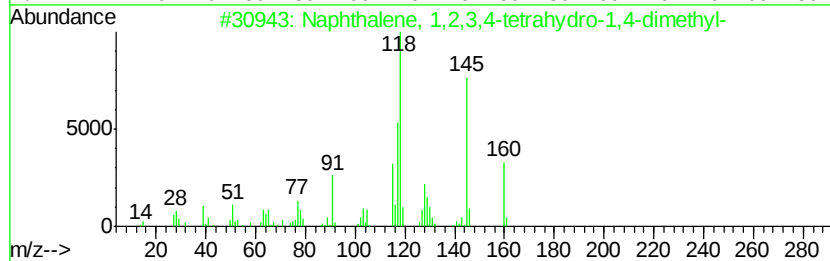
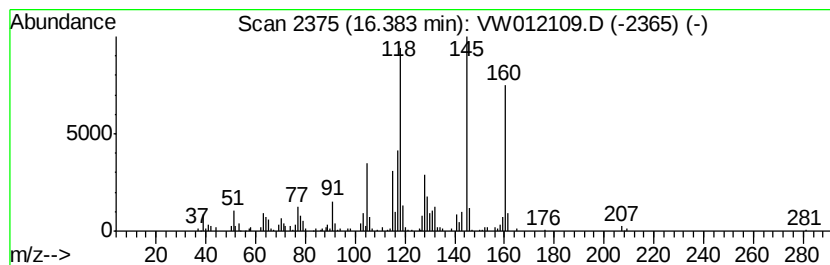
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	10.37 ug/l	129721	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro...	160	C12H16	004175-54-6	95
2		Naphthalene, 1,2,3,4-tetrahydro...	160	C12H16	013065-07-1	81
3		Naphthalene, 1,2,3,4-tetrahydro...	160	C12H16	007524-63-2	76
4		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	64
5		Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW082219\
 Data File : VW012109.D
 Acq On : 22 Aug 2019 14:34
 Operator : SY/VA
 Sample : K4086-01
 Misc : 4.01G/5ML/MSVOA_W/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 NB-08-082019

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.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
Benzene, 2-etheny...	14.81	12.7	ug/l	158235	4	13.56	625417	50.0
Naphthalene, 1,2,...	14.96	6.3	ug/l	78973	4	13.56	625417	50.0
Benzene, (1,2-dim...	15.07	6.2	ug/l	77346	4	13.56	625417	50.0
1H-Indene, 2,3-di...	15.18	6.5	ug/l	80761	4	13.56	625417	50.0
Disilane, hexamet...	15.50	15.6	ug/l	194843	4	13.56	625417	50.0
1H-Indene, 2,3-di...	15.66	7.0	ug/l	87297	4	13.56	625417	50.0
1H-Indene, 2,3-di...	15.83	6.0	ug/l	74441	4	13.56	625417	50.0
Naphthalene, 1,2,...	15.86	13.1	ug/l	163702	4	13.56	625417	50.0
Naphthalene, 1,2,...	16.38	10.4	ug/l	129721	4	13.56	625417	50.0