

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW111318\
 Data File : VW006869.D
 Acq On : 13 Nov 2018 21:34
 Operator : SY/AP
 Sample : J5860-21
 Misc : 5.47G/5ML/MSVOA W/SOIL
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EP2-SB-110(14-16)

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\82W111218S.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	4.922	483	495	510	rVB	26913	86101	2.32%	0.168%
2	7.891	972	982	986	rBV	181699	422790	11.39%	0.826%
3	7.952	986	992	1004	rVB	486697	1078546	29.06%	2.108%
4	8.311	1041	1051	1062	rBV	158848	345079	9.30%	0.674%
5	8.848	1129	1139	1150	rBV	603854	1188837	32.03%	2.323%
6	10.329	1374	1382	1389	rBV	772752	1377986	37.13%	2.693%
7	10.756	1440	1452	1463	rVB2	59938	235590	6.35%	0.460%
8	11.414	1553	1560	1571	rVB6	16061	43291	1.17%	0.085%
9	11.634	1587	1596	1604	rBV	907517	1542381	41.56%	3.014%
10	11.738	1605	1613	1622	rVB	35550	64466	1.74%	0.126%
11	11.847	1622	1631	1641	rBV	63800	144268	3.89%	0.282%
12	12.140	1673	1679	1685	rBV4	20273	42338	1.14%	0.083%
13	12.335	1702	1711	1714	rBV5	15743	37475	1.01%	0.073%
14	12.622	1747	1758	1767	rVB	694415	1310386	35.31%	2.561%
15	12.762	1776	1781	1784	rBV3	22132	43884	1.18%	0.086%
16	12.811	1784	1789	1793	rVB	77516	126762	3.42%	0.248%
17	12.871	1793	1799	1801	rBV	115944	174168	4.69%	0.340%
18	12.902	1801	1804	1808	rVV	169759	284490	7.67%	0.556%
19	12.951	1808	1812	1818	rVB	303301	486294	13.10%	0.950%
20	13.109	1829	1838	1845	rBV	173848	381159	10.27%	0.745%
21	13.170	1845	1848	1850	rBV	36522	46943	1.26%	0.092%
22	13.201	1850	1853	1856	rVB	43314	52137	1.40%	0.102%
23	13.256	1856	1862	1867	rBV	1465235	2194390	59.13%	4.288%
24	13.304	1867	1870	1874	rVB2	55272	78844	2.12%	0.154%
25	13.390	1874	1884	1885	rBV4	51051	121718	3.28%	0.238%
26	13.414	1885	1888	1891	rBV	61590	69890	1.88%	0.137%
27	13.451	1891	1894	1897	rVB2	55500	72129	1.94%	0.141%
28	13.493	1897	1901	1904	rBV2	78433	123585	3.33%	0.242%
29	13.560	1907	1912	1916	rBV	1049512	1867263	50.31%	3.649%
30	13.725	1934	1939	1941	rBV	267628	406726	10.96%	0.795%
31	13.762	1941	1945	1948	rVV2	890680	1536480	41.40%	3.003%
32	13.798	1948	1951	1961	rVB2	778007	1445118	38.94%	2.824%
33	13.883	1961	1965	1970	rBV2	139116	247534	6.67%	0.484%
34	13.957	1972	1977	1981	rVB	236759	349986	9.43%	0.684%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Title : SW846 8260

35	14.018	1981	1987	1989	rBV	574303	856613	23.08%	1.674%
36	14.048	1989	1992	1996	rVB	447124	618644	16.67%	1.209%
37	14.103	1996	2001	2007	rBV	1109002	1663766	44.83%	3.251%
38	14.164	2008	2011	2014	rVB	105432	126034	3.40%	0.246%
39	14.213	2014	2019	2024	rVB2	733991	1169887	31.52%	2.286%
40	14.268	2024	2028	2035	rVB6	91884	220030	5.93%	0.430%
41	14.347	2035	2041	2046	rBV3	360501	655338	17.66%	1.281%
42	14.426	2046	2054	2057	rBV	815731	1449562	39.06%	2.833%
43	14.463	2057	2060	2064	rVB	1056944	1380265	37.19%	2.697%
44	14.505	2064	2067	2071	rBV2	649266	914851	24.65%	1.788%
45	14.548	2071	2074	2081	rVB	366683	534293	14.40%	1.044%
46	14.609	2081	2084	2086	rBV	46920	55020	1.48%	0.108%
47	14.652	2086	2091	2094	rBV	273493	387999	10.45%	0.758%
48	14.694	2094	2098	2102	rBV	826933	1358172	36.60%	2.654%
49	14.737	2102	2105	2110	rVB2	487202	738067	19.89%	1.442%
50	14.816	2110	2118	2123	rBV3	1509365	3711298	100.00%	7.252%
51	14.969	2139	2143	2150	rVB3	163345	325197	8.76%	0.635%
52	15.072	2150	2160	2164	rBV2	980530	2308222	62.19%	4.511%
53	15.188	2172	2179	2187	rVB2	796915	1940324	52.28%	3.792%
54	15.261	2187	2191	2198	rVB5	177699	323981	8.73%	0.633%
55	15.341	2198	2204	2206	rBV3	698500	1256119	33.85%	2.455%
56	15.371	2206	2209	2215	rVB	892074	1465768	39.49%	2.864%
57	15.481	2221	2227	2231	rBV2	432490	846912	22.82%	1.655%
58	15.530	2231	2235	2238	rVV2	212629	361657	9.74%	0.707%
59	15.670	2250	2258	2264	rBV3	655937	1698537	45.77%	3.319%
60	15.956	2302	2305	2310	rVB2	218104	357486	9.63%	0.699%
61	16.029	2310	2317	2324	rBV3	591880	1642563	44.26%	3.210%
62	16.176	2334	2341	2346	rBV8	203774	567862	15.30%	1.110%
63	16.237	2347	2351	2356	rVV5	234799	456927	12.31%	0.893%
64	16.298	2357	2361	2367	rVB	465687	771188	20.78%	1.507%
65	16.450	2381	2386	2391	rBV2	527549	1057487	28.49%	2.066%
66	16.511	2392	2396	2402	rVB2	300991	514716	13.87%	1.006%
67	16.651	2412	2419	2432	rVB4	402128	1409514	37.98%	2.754%

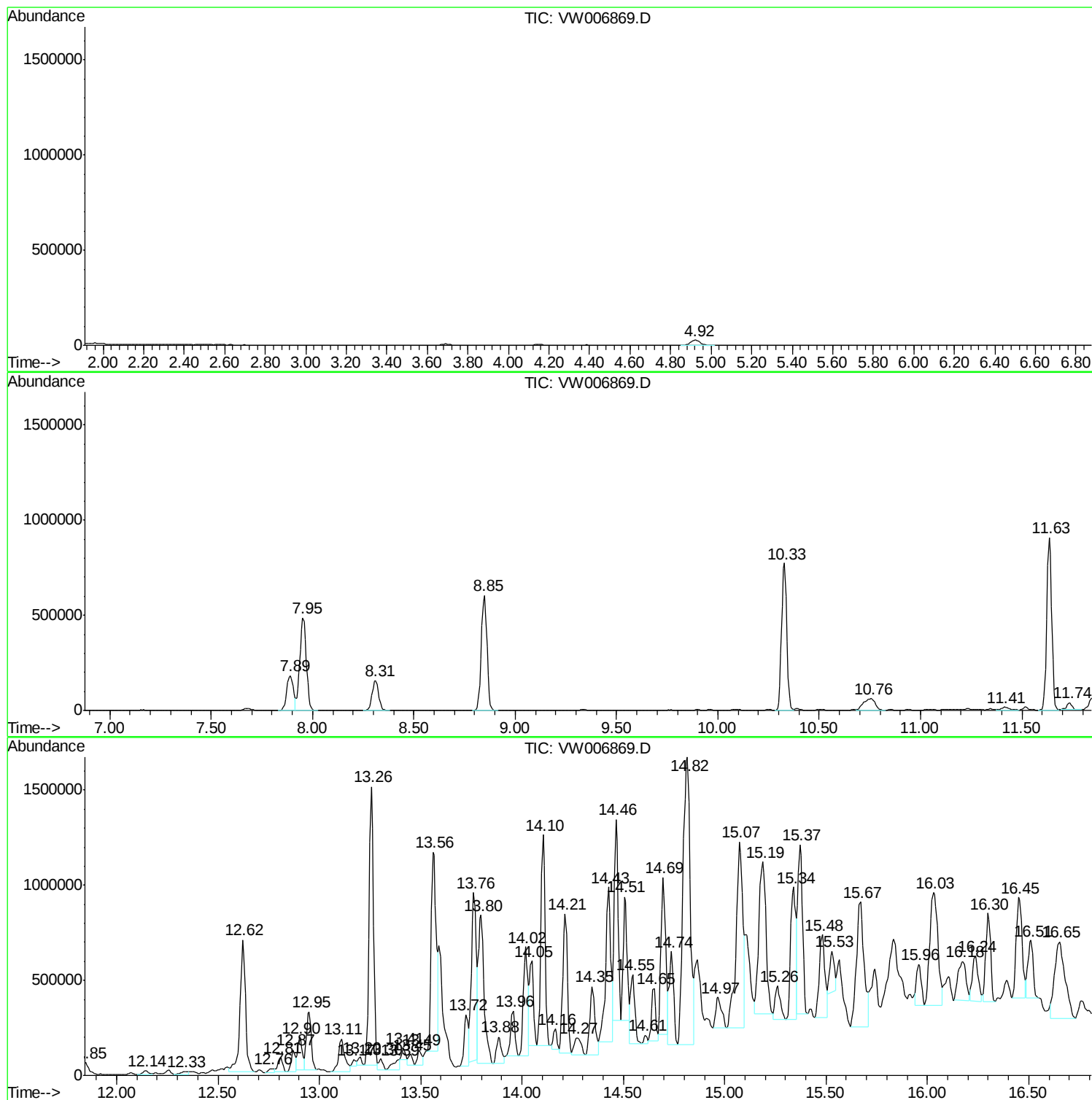
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Instrument :
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ClientSampleId :
EP2-SB-110(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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



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ClientSampleId :
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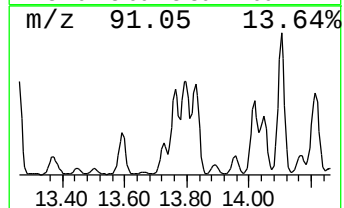
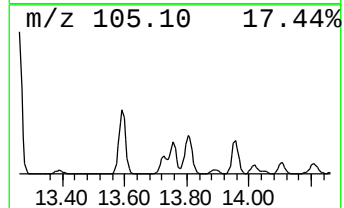
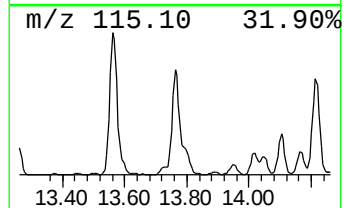
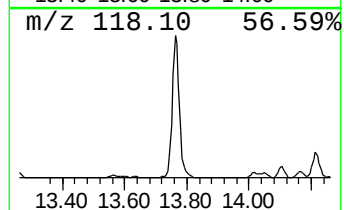
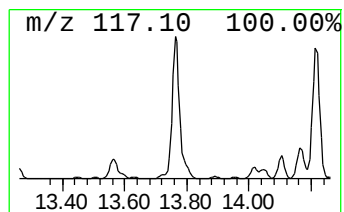
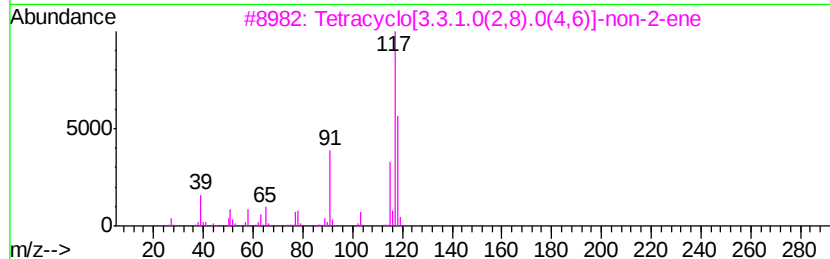
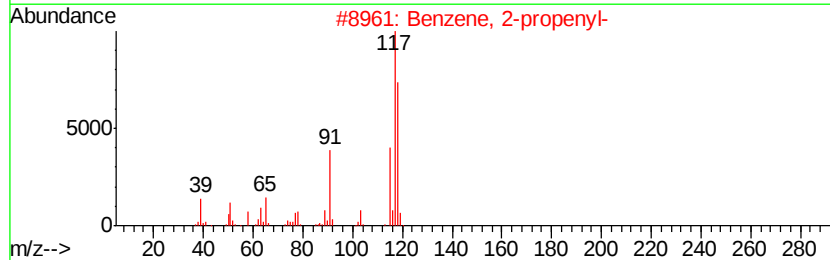
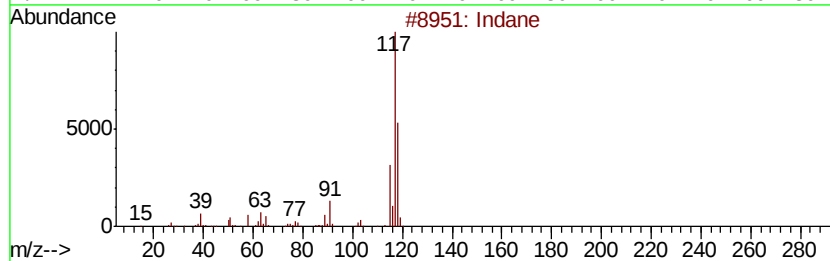
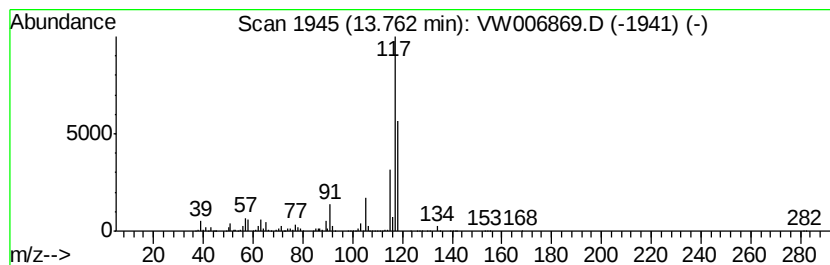
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 1 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	41.14 ug/l	1536480	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	93
2	Benzene, 2-propenyl-		118	C9H10	000300-57-2	81
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	72
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	72
5	1,4:5,8-Dimethanobiphenylene, 1,	...	184	C14H16	001624-13-1	64



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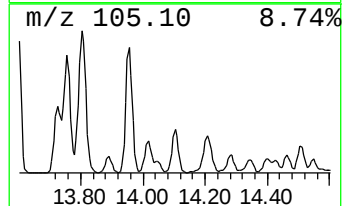
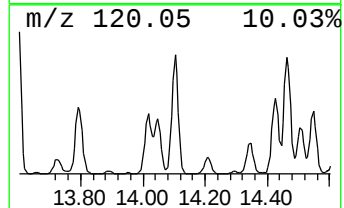
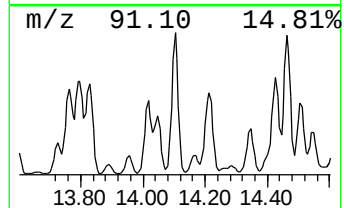
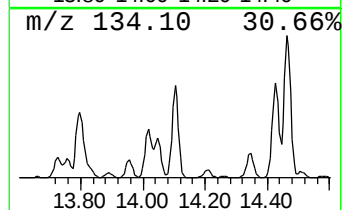
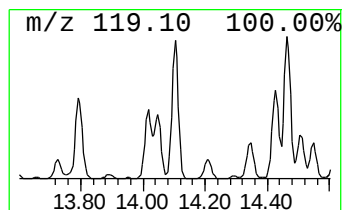
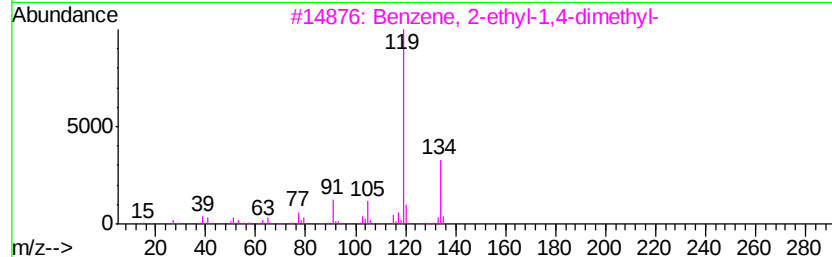
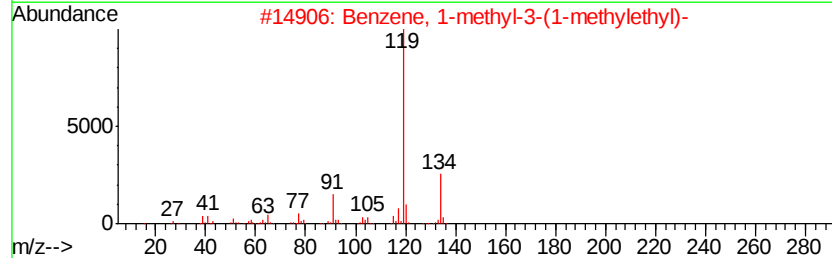
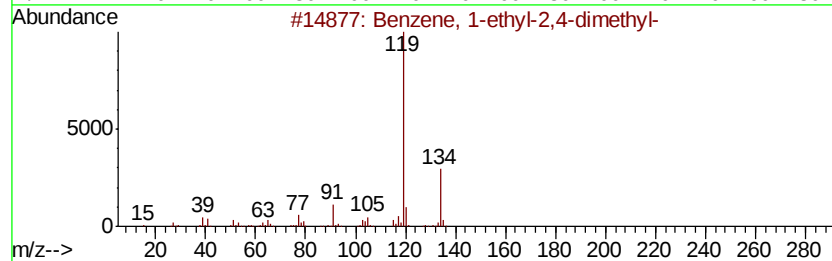
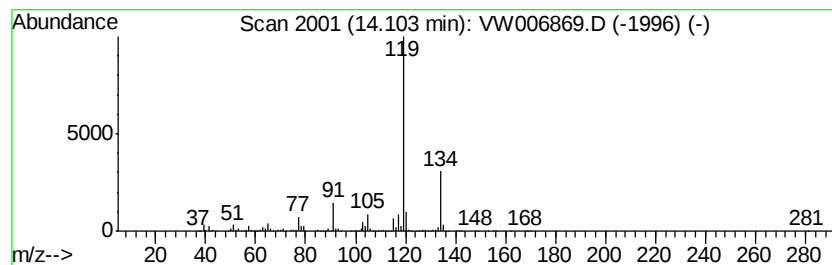
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	44.55 ug/l	1663770	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



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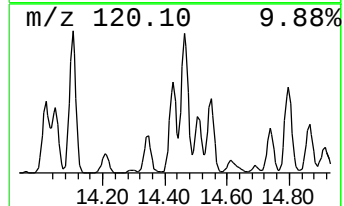
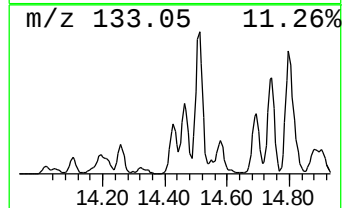
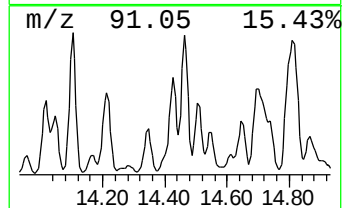
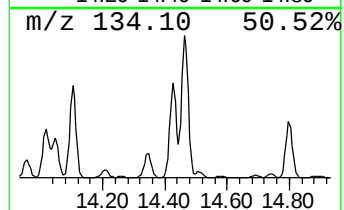
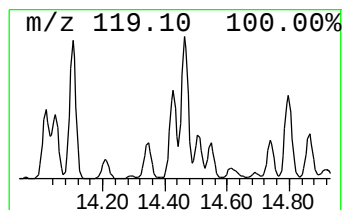
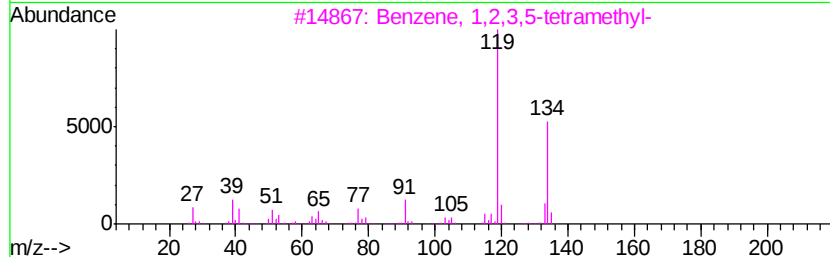
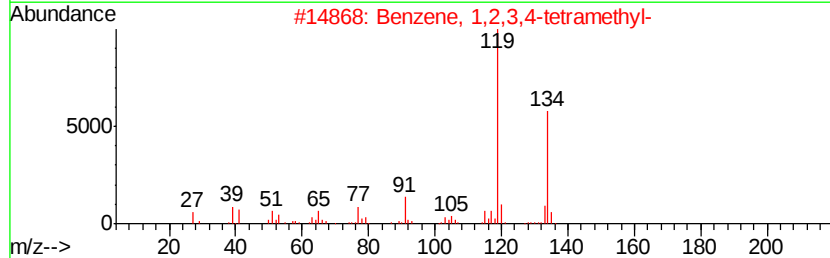
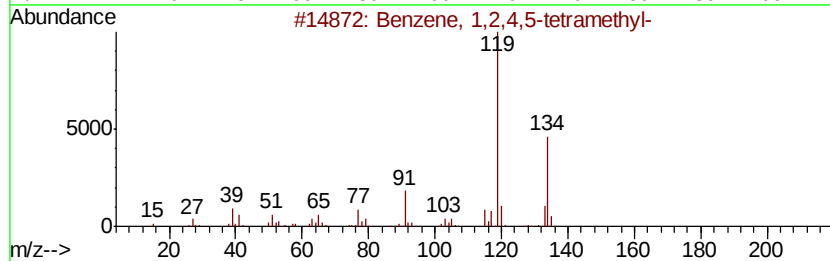
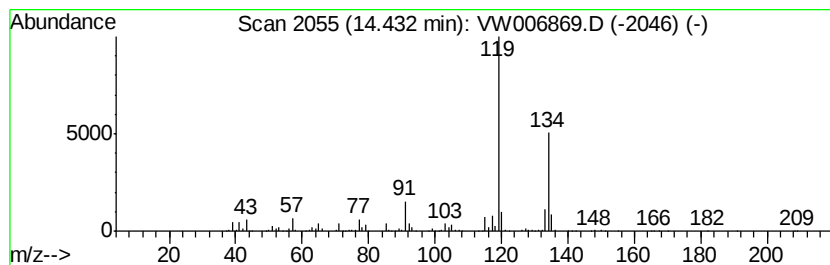
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	38.82 ug/l	1449560	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91



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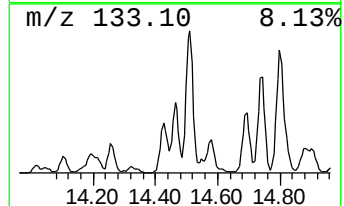
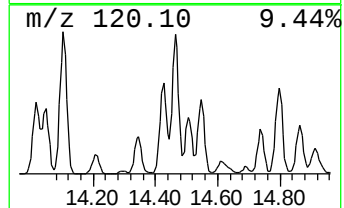
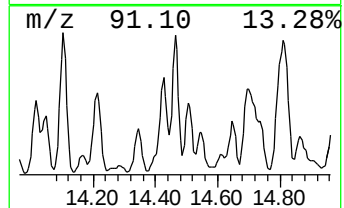
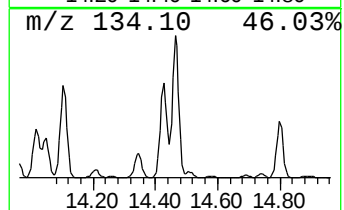
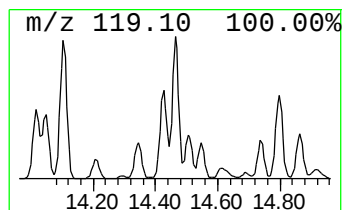
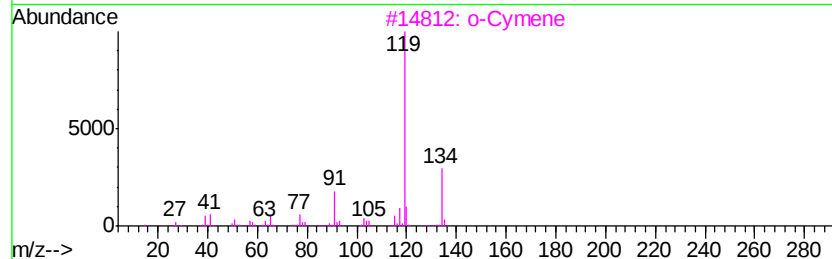
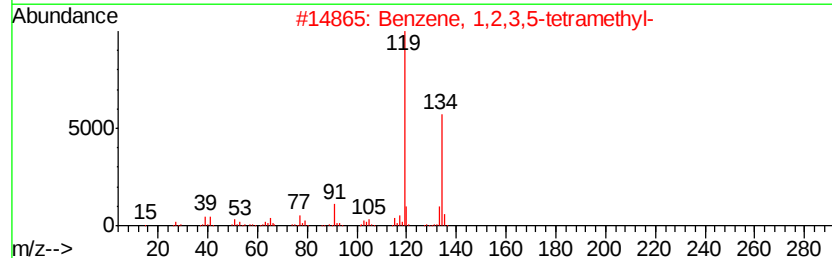
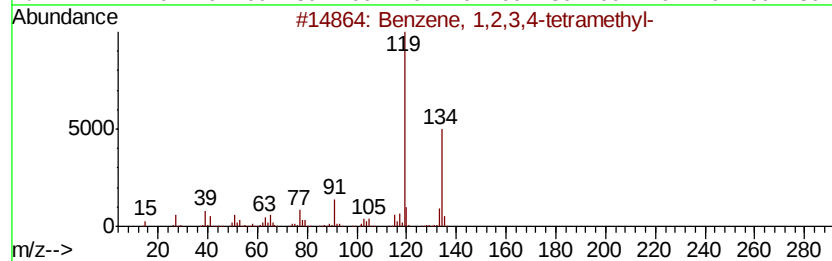
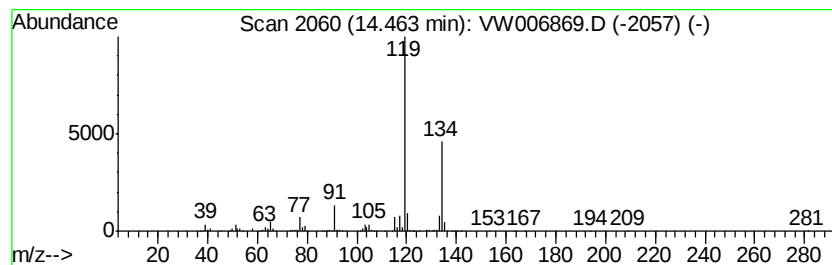
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	36.96 ug/l	1380270	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		o-Cymene	134	C10H14	000527-84-4	94
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

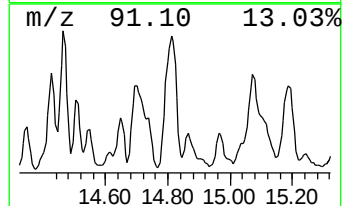
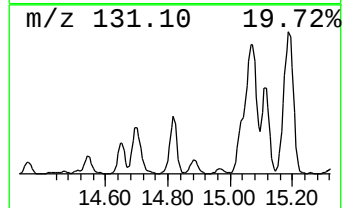
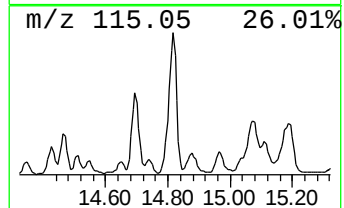
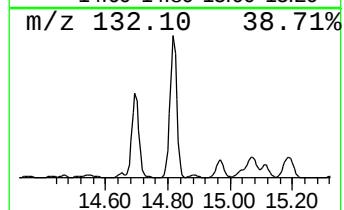
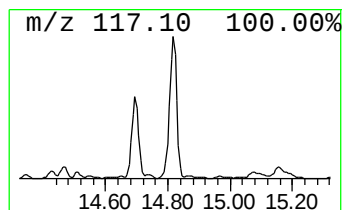
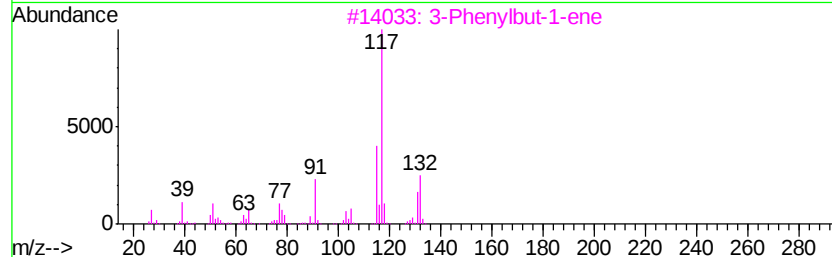
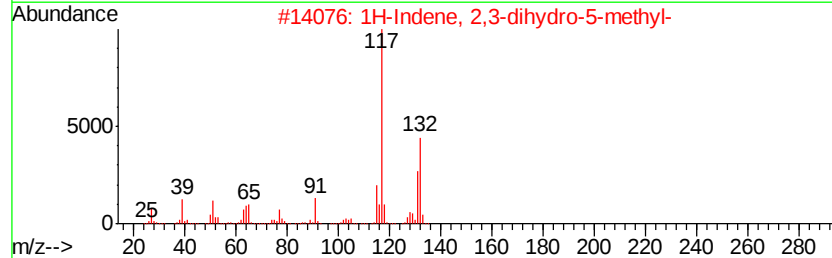
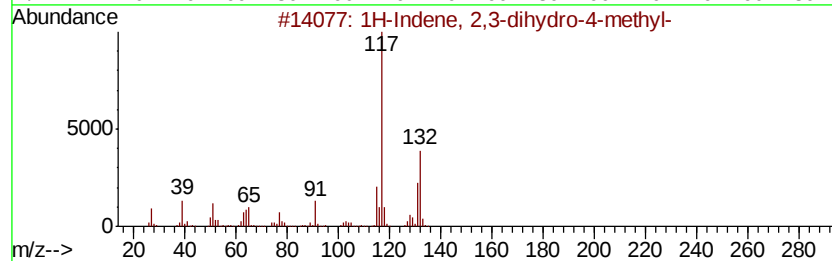
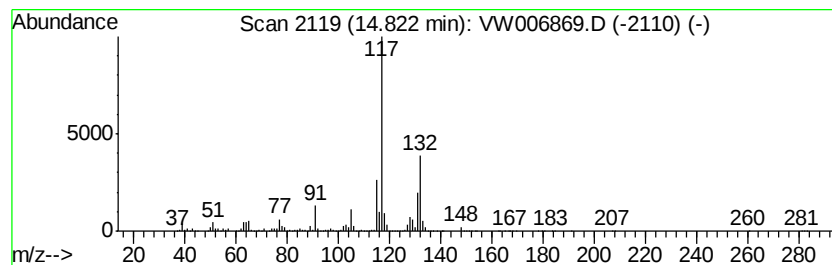
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	99.38 ug/l	3711300	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

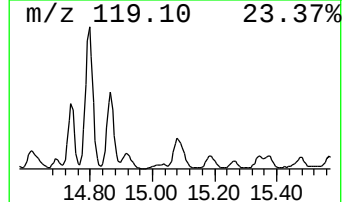
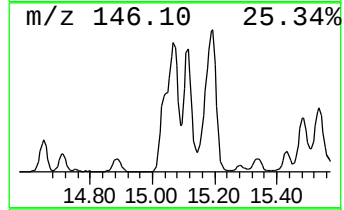
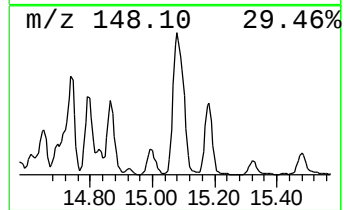
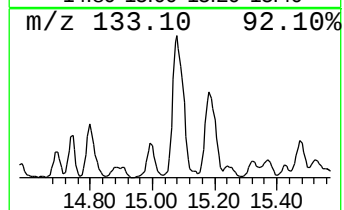
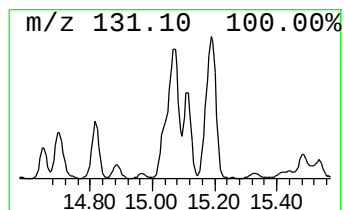
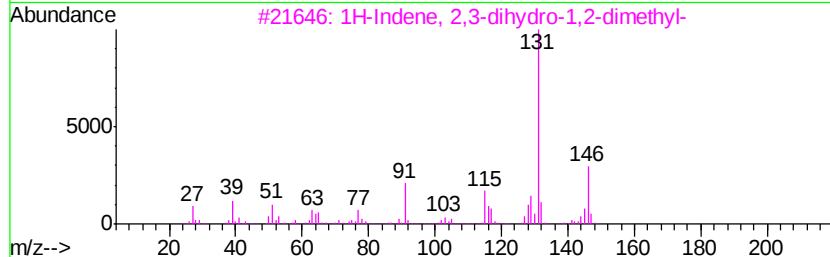
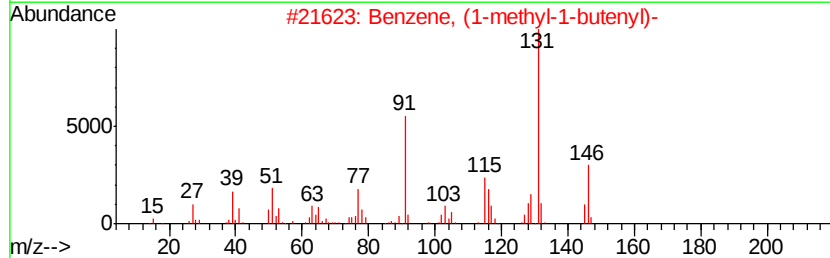
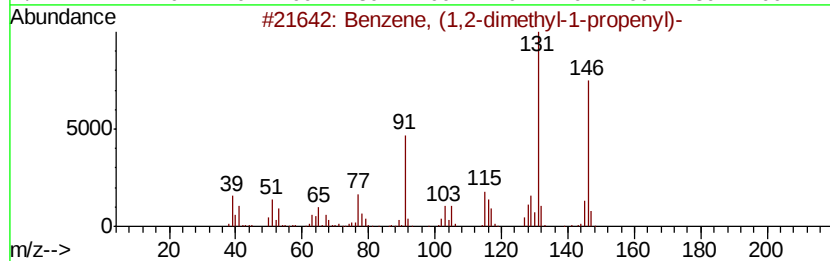
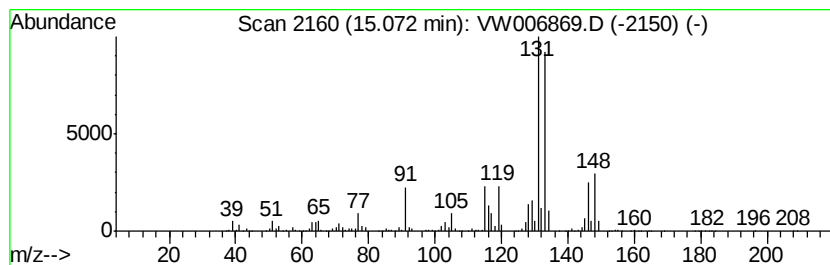
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	61.81 ug/l	2308220	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
2		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
3		1H-Indene, 2,3-dihydro-1,2-dimethyl-	146	C11H14	017057-82-8	89
4		1H-Indene, 2,3-dihydro-1,6-dimethyl-	146	C11H14	017059-48-2	86
5		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

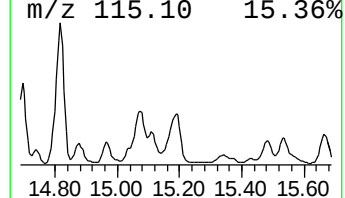
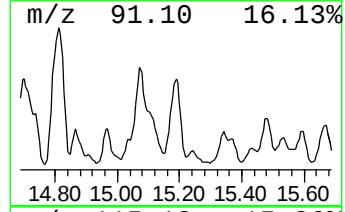
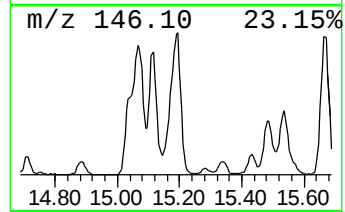
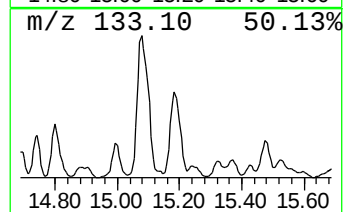
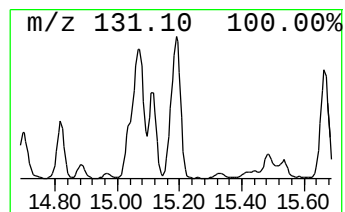
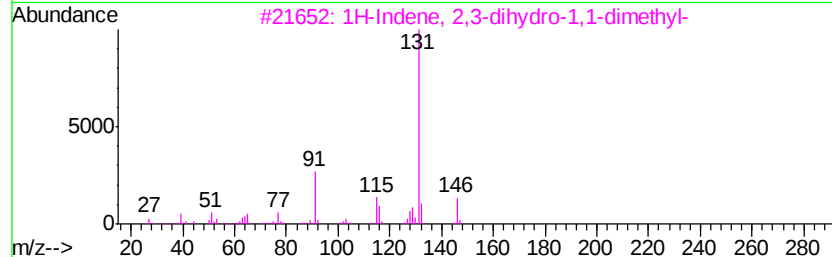
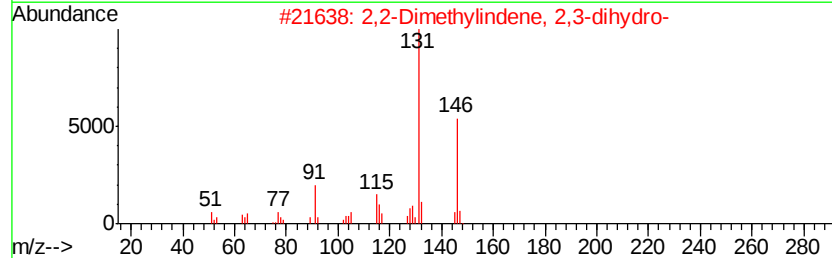
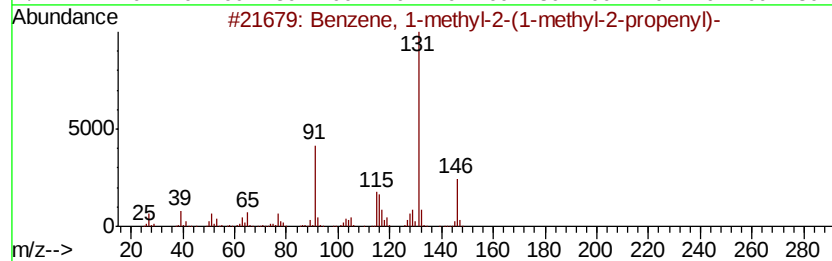
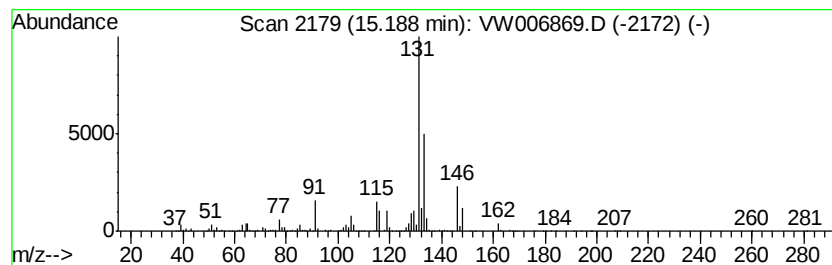
Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-2-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	51.96 ug/l	1940320	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-2-(1-methyl-2-...	146 C11H14	097664-19-2 70
2		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 64
3		1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9 64
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 64
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

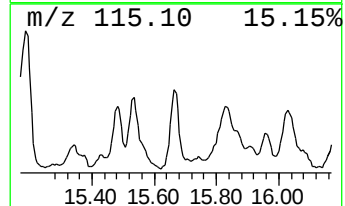
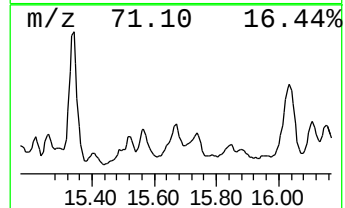
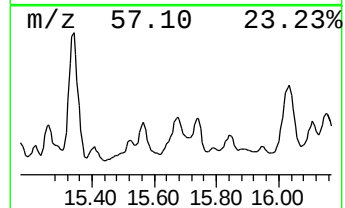
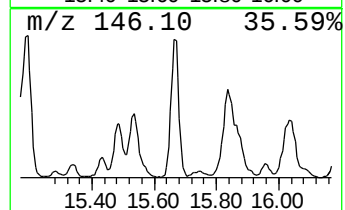
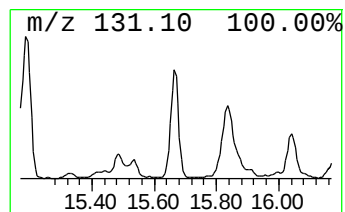
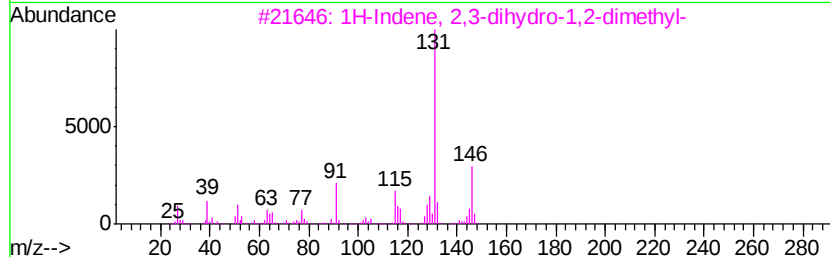
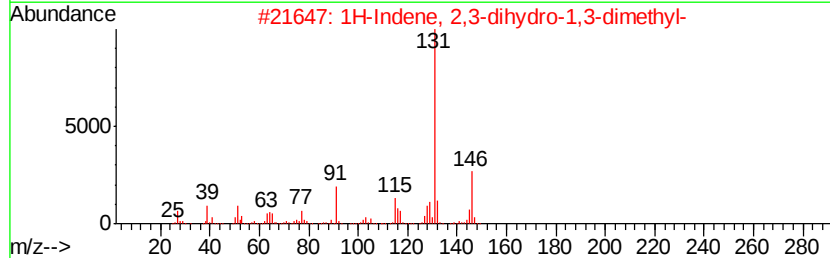
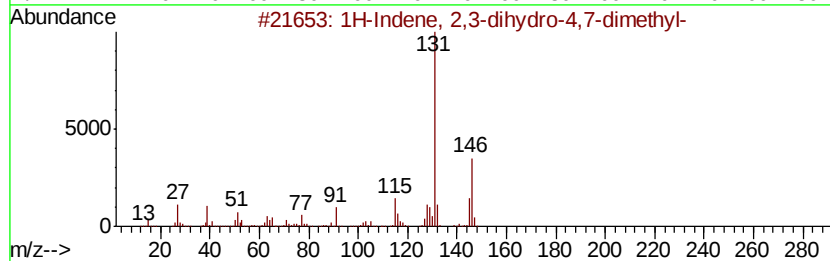
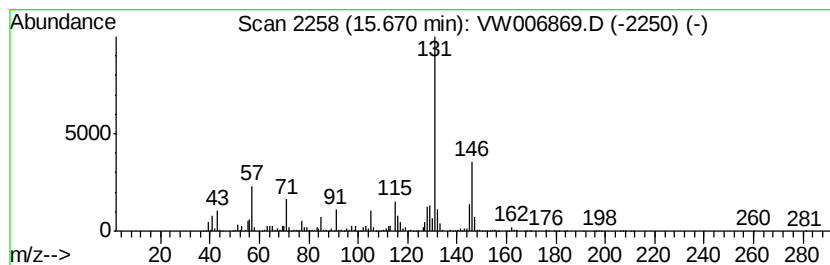
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	45.48 ug/l	1698540	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	97
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
3		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

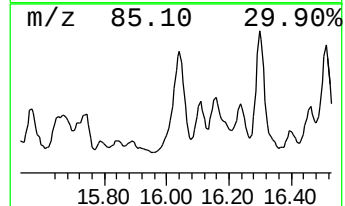
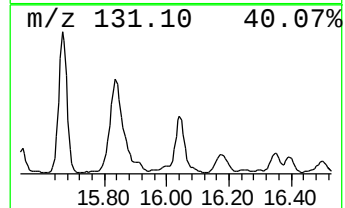
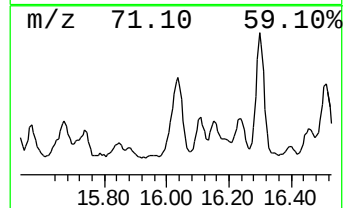
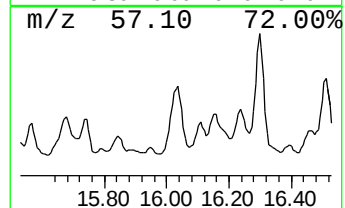
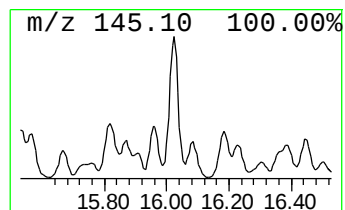
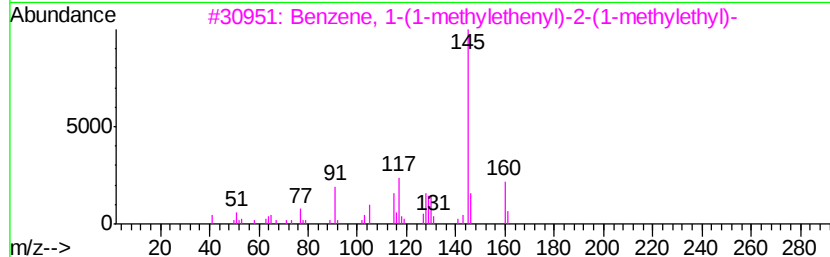
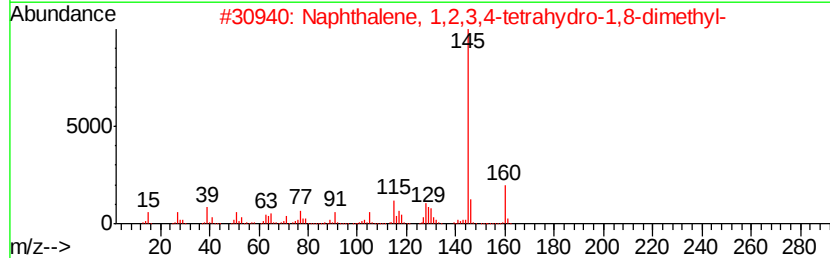
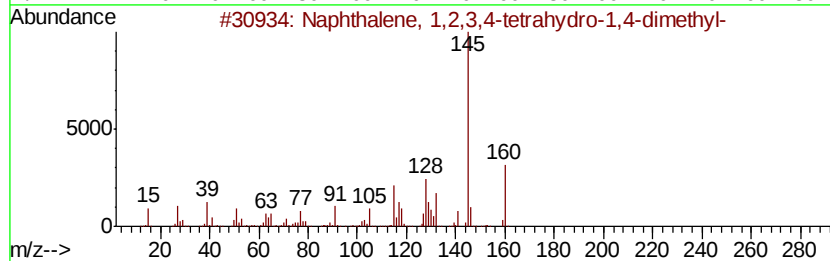
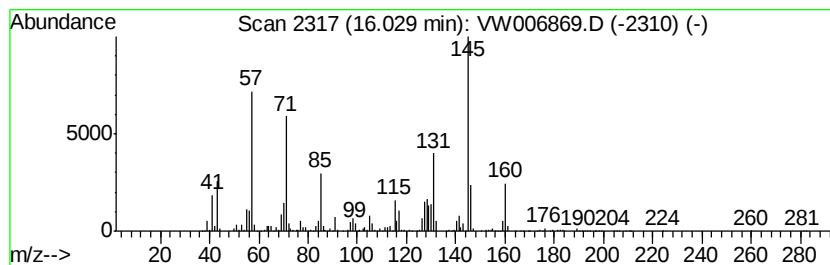
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	43.98 ug/l	1642560	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	50
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	50
3		Benzene, 1-(1-methylethenyl)-2-(...	160	C12H16	005557-93-7	49
4		1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	49
5		1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

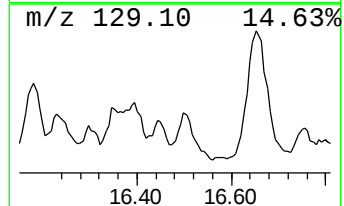
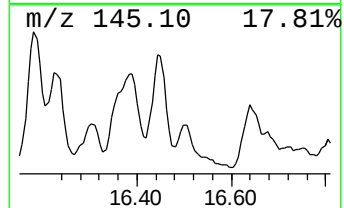
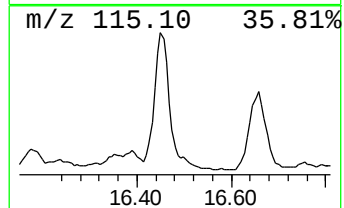
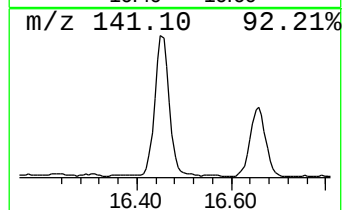
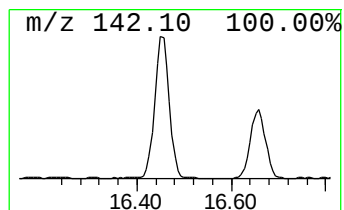
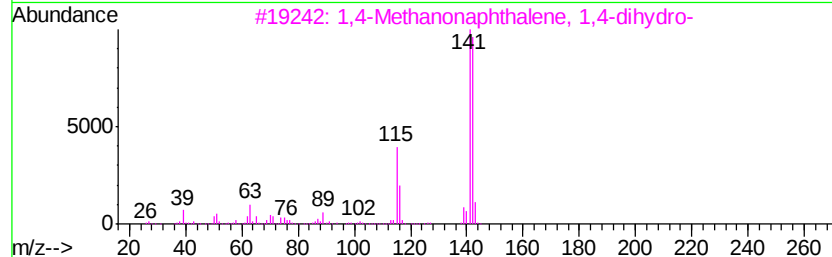
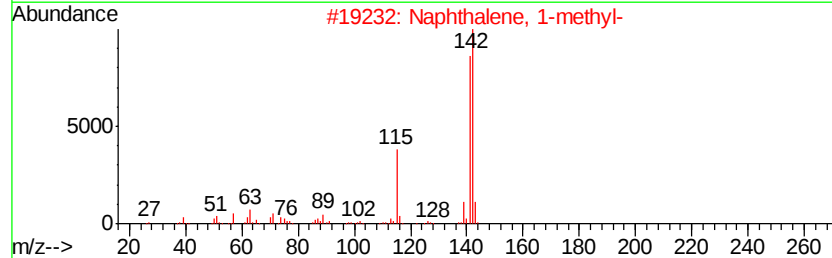
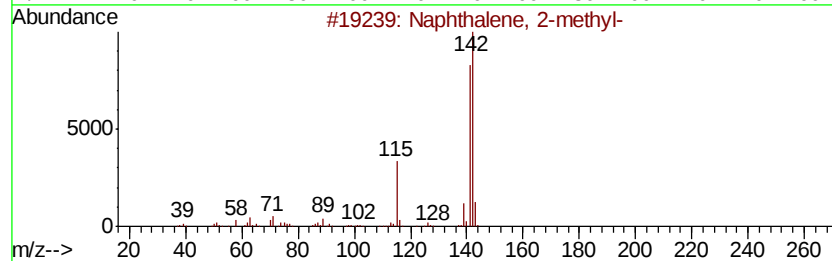
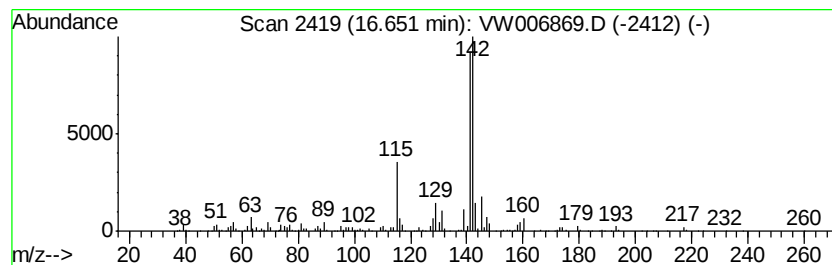
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	37.74 ug/l	1409510	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	81
4		Benzocycloheptatriene	142	C11H10	000264-09-5	74
5		Spiro(9-methylenetricyclo[6.2.1....	184	C13H12O	033929-47-4	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW111318\
Data File : VW006869.D
Acq On : 13 Nov 2018 21:34
Operator : SY/AP
Sample : J5860-21
Misc : 5.47G/5ML/MSVOA_W/SOIL
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-110(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.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
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Benzene, 1-ethyl-...	14.10	44.5	ug/l	1663770	4	13.57	1867260	50.0
Benzene, 1,2,4,5-...	14.43	38.8	ug/l	1449560	4	13.57	1867260	50.0
Benzene, 1,2,3,4-...	14.46	37.0	ug/l	1380270	4	13.57	1867260	50.0
1H-Indene, 2,3-di...	14.82	99.4	ug/l	3711300	4	13.57	1867260	50.0
Benzene, (1,2-dim...	15.07	61.8	ug/l	2308220	4	13.57	1867260	50.0
Benzene, 1-methyl...	15.19	52.0	ug/l	1940320	4	13.57	1867260	50.0
1H-Indene, 2,3-di...	15.67	45.5	ug/l	1698540	4	13.57	1867260	50.0
Naphthalene, 1,2,...	16.03	44.0	ug/l	1642560	4	13.57	1867260	50.0
Naphthalene, 1-me...	16.65	37.7	ug/l	1409510	4	13.57	1867260	50.0