

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD082119\
 Data File : VD063722.D
 Acq On : 21 Aug 2019 19:28
 Operator : JC/SY
 Sample : K4418-04
 Misc : 5.00µl/5ml/MSVOA D/SOIL
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CFVP-228L

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_D\METHOD\82D080919S.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	1.779	30	36	37	rBV3	9532	27672	1.37%	0.064%
2	3.207	178	181	188	rVB3	20988	56390	2.79%	0.131%
3	6.918	553	558	565	rBV	156263	442072	21.89%	1.030%
4	7.046	565	571	584	rVB	150111	422452	20.92%	0.984%
5	7.449	606	612	619	rBV	254444	652115	32.29%	1.520%
6	8.217	684	690	701	rBV	195989	498343	24.67%	1.161%
7	8.857	750	755	763	rVB5	16939	40210	1.99%	0.094%
8	10.413	907	913	919	rBV2	266943	602536	29.83%	1.404%
9	10.511	919	923	931	rVB	94787	220823	10.93%	0.515%
10	12.185	1089	1093	1097	rVB	21224	42842	2.12%	0.100%
11	12.313	1100	1106	1108	rBV2	22234	35971	1.78%	0.084%
12	12.372	1108	1112	1118	rVB	240638	408510	20.22%	0.952%
13	12.529	1124	1128	1134	rVB	70694	116501	5.77%	0.271%
14	12.667	1134	1142	1147	rVV	373199	691115	34.22%	1.610%
15	12.746	1147	1150	1155	rVB	47979	81672	4.04%	0.190%
16	13.002	1171	1176	1178	rBV3	20091	40790	2.02%	0.095%
17	13.051	1178	1181	1184	rVV	406317	570552	28.25%	1.329%
18	13.090	1184	1185	1188	rVB2	17388	20620	1.02%	0.048%
19	13.218	1192	1198	1201	rBV2	28538	45726	2.26%	0.107%
20	13.307	1204	1207	1213	rVB7	30859	69525	3.44%	0.162%
21	13.405	1213	1217	1220	rVB2	61133	112520	5.57%	0.262%
22	13.484	1220	1225	1227	rBV4	23919	47774	2.37%	0.111%
23	13.553	1227	1232	1237	rBV3	259060	426049	21.09%	0.993%
24	13.661	1240	1243	1246	rVB	48775	68951	3.41%	0.161%
25	13.710	1246	1248	1251	rVB3	28270	33695	1.67%	0.079%
26	13.779	1251	1255	1257	rBV	49561	71069	3.52%	0.166%
27	13.848	1259	1262	1264	rBV	522816	640702	31.72%	1.493%
28	13.878	1264	1265	1267	rVB	124374	115654	5.73%	0.269%
29	13.937	1267	1271	1273	rBV	348386	558694	27.66%	1.302%
30	13.966	1273	1274	1278	rVB2	190970	212247	10.51%	0.495%
31	14.084	1282	1286	1288	rBV	310326	400145	19.81%	0.932%
32	14.114	1288	1289	1292	rVB3	33048	39396	1.95%	0.092%
33	14.163	1292	1294	1296	rBV2	17954	32478	1.61%	0.076%
34	14.203	1296	1298	1299	rBV	51502	56363	2.79%	0.131%

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35	14.242	1299	1302	1308	rVB	1367260	2019827	100.00%	4.706%
36	14.331	1308	1311	1312	rBV	21786	37703	1.87%	0.088%
37	14.370	1312	1315	1319	rVB4	67399	105170	5.21%	0.245%
38	14.449	1319	1323	1326	rBV2	198568	490257	24.27%	1.142%
39	14.488	1326	1327	1328	rBV	63976	56482	2.80%	0.132%
40	14.518	1328	1330	1332	rVB	197134	203742	10.09%	0.475%
41	14.557	1332	1334	1339	rVB	713604	903924	44.75%	2.106%
42	14.646	1341	1343	1346	rVB	108821	136379	6.75%	0.318%
43	14.705	1346	1349	1354	rBV2	465540	904839	44.80%	2.108%
44	14.774	1354	1356	1359	rBV	393662	600325	29.72%	1.399%
45	14.882	1363	1367	1369	rBV	391774	580821	28.76%	1.353%
46	14.980	1374	1377	1378	rVV	339456	489687	24.24%	1.141%
47	15.000	1378	1379	1383	rVV	285404	533549	26.42%	1.243%
48	15.059	1383	1385	1387	rVV	489530	648583	32.11%	1.511%
49	15.089	1387	1388	1390	rVV	150156	187740	9.29%	0.437%
50	15.138	1390	1393	1398	rVV2	521761	947520	46.91%	2.208%
51	15.226	1398	1402	1404	rVV4	122081	306552	15.18%	0.714%
52	15.266	1404	1406	1409	rVV2	226046	477977	23.66%	1.114%
53	15.345	1409	1414	1416	rVV3	382921	936061	46.34%	2.181%
54	15.384	1416	1418	1420	rVV	566184	819052	40.55%	1.908%
55	15.433	1420	1423	1425	rVV2	276200	546652	27.06%	1.274%
56	15.473	1425	1427	1430	rVV3	156610	331124	16.39%	0.772%
57	15.571	1430	1437	1441	rVV2	500204	1253991	62.08%	2.922%
58	15.650	1441	1445	1446	rVV	658879	1213854	60.10%	2.828%
59	15.669	1446	1447	1452	rVV	1137381	1811949	89.71%	4.222%
60	15.748	1452	1455	1457	rVV2	360769	679566	33.64%	1.583%
61	15.797	1457	1460	1463	rVV	1076960	1651468	81.76%	3.848%
62	15.896	1466	1470	1472	rVV	325314	947195	46.89%	2.207%
63	15.935	1472	1474	1476	rVV	345898	597030	29.56%	1.391%
64	15.994	1476	1480	1482	rVV	493085	1150797	56.98%	2.682%
65	16.053	1484	1486	1487	rVV	285059	449913	22.27%	1.048%
66	16.083	1487	1489	1491	rVV2	276070	523105	25.90%	1.219%
67	16.142	1491	1495	1497	rVV2	440036	1093383	54.13%	2.548%
68	16.171	1497	1498	1501	rVV	411058	695099	34.41%	1.620%
69	16.231	1501	1504	1509	rVV3	587975	1444336	71.51%	3.365%
70	16.319	1509	1513	1515	rVV2	485572	1025605	50.78%	2.390%
71	16.359	1515	1517	1521	rVV	350654	809893	40.10%	1.887%

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72	16.418	1521	1523	1526	rVV3	212048	497998	24.66%	1.160%
73	16.506	1526	1532	1535	rVV	916099	1890743	93.61%	4.406%
74	16.575	1537	1539	1542	rVV3	253001	536508	26.56%	1.250%
75	16.624	1542	1544	1551	rVV	253743	864082	42.78%	2.013%
76	16.733	1551	1555	1558	rVV	513777	976659	48.35%	2.276%
77	16.782	1558	1560	1564	rVV5	117304	326663	16.17%	0.761%
78	16.880	1564	1570	1572	rVV3	280755	755976	37.43%	1.762%
79	16.920	1572	1574	1579	rVV3	247187	522888	25.89%	1.218%
80	16.989	1579	1581	1583	rVV2	82679	149033	7.38%	0.347%
81	17.057	1583	1588	1591	rVV2	173386	428758	21.23%	0.999%
82	17.107	1591	1593	1595	rVV3	67140	110894	5.49%	0.258%
83	17.166	1595	1599	1605	rVV7	51013	158526	7.85%	0.369%
84	17.254	1605	1608	1611	rVB4	46580	84029	4.16%	0.196%
85	17.382	1619	1621	1624	rVB4	16198	26776	1.33%	0.062%
86	17.461	1624	1629	1633	rVV7	27095	73261	3.63%	0.171%

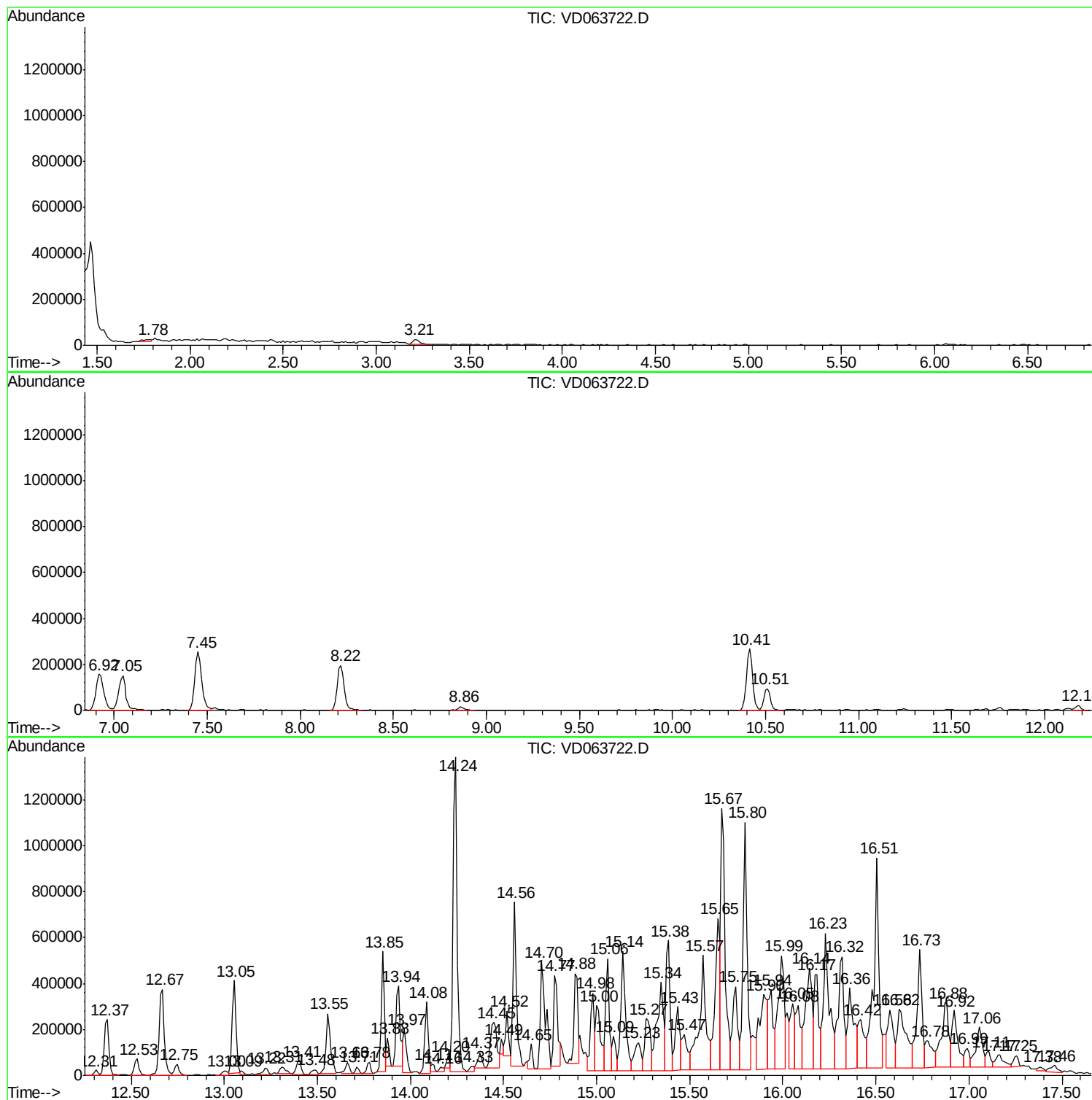
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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



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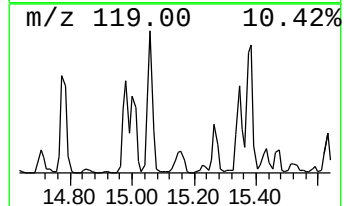
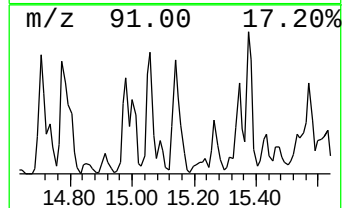
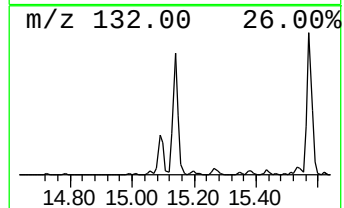
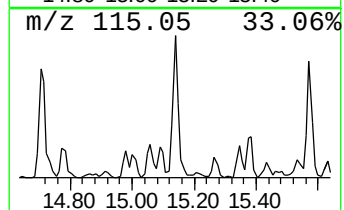
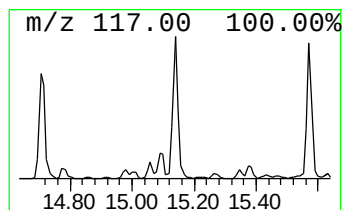
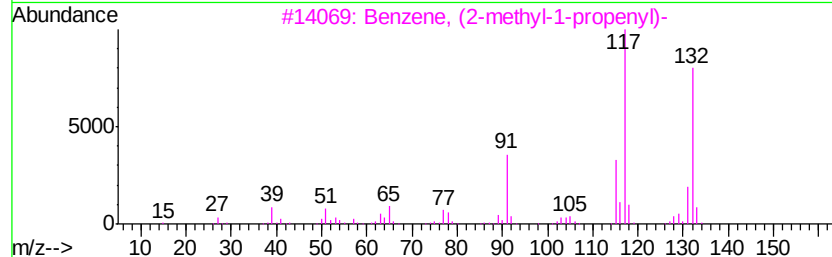
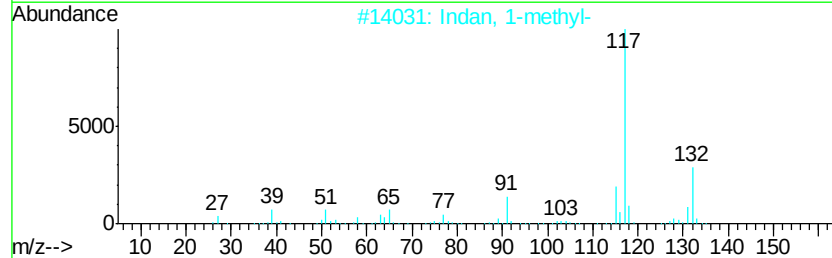
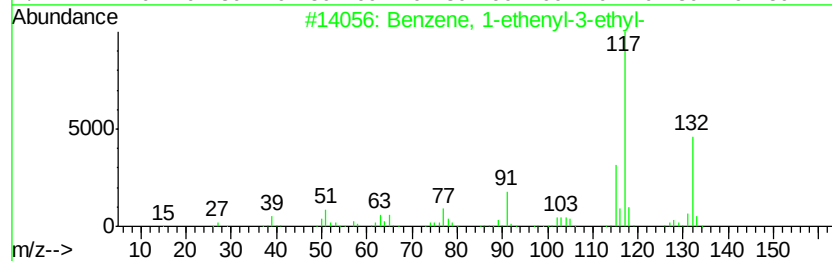
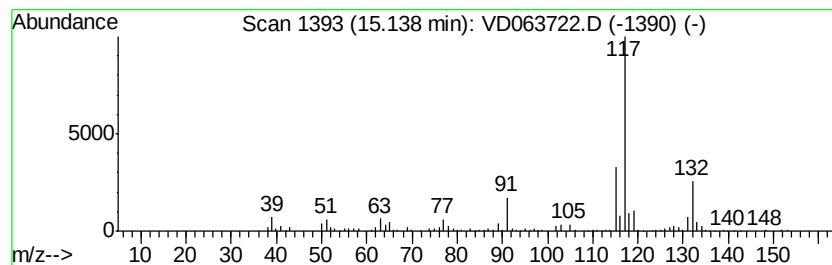
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	232.53 ug/l	947520	1,4-Dichlorobenzene-d4	14.52

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	87
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	72
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	64
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	60



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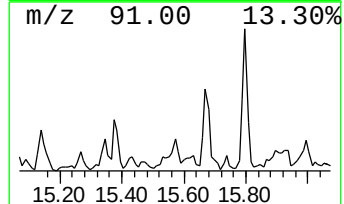
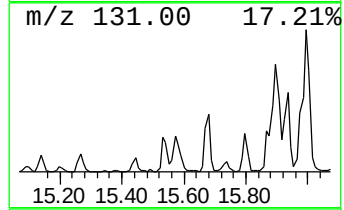
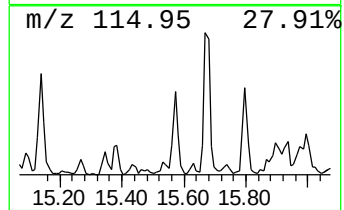
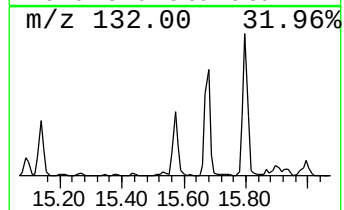
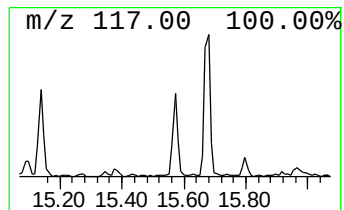
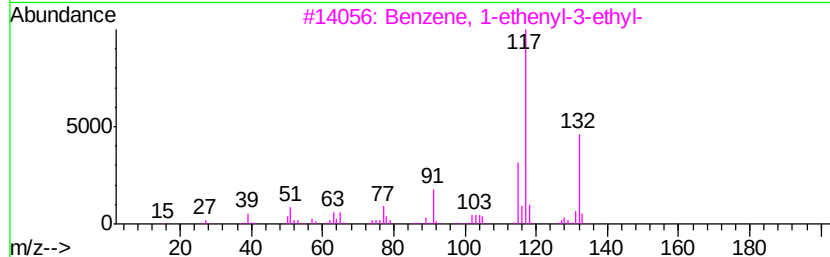
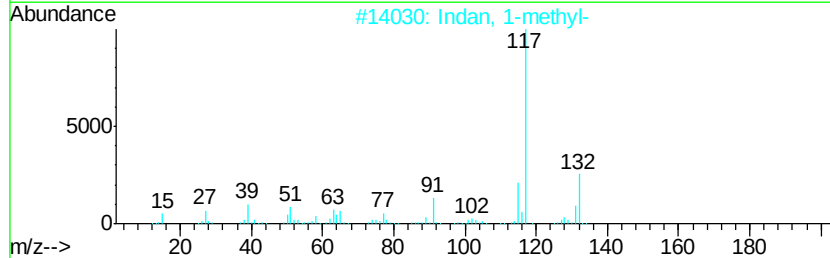
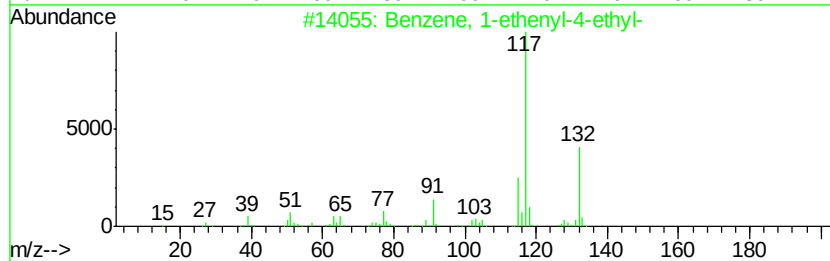
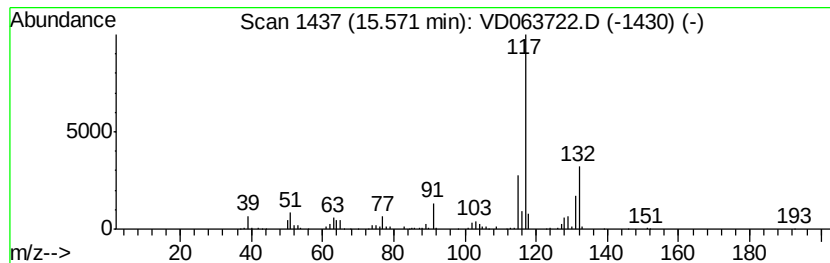
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	307.74 ug/l	1253990	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86



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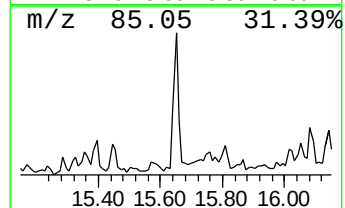
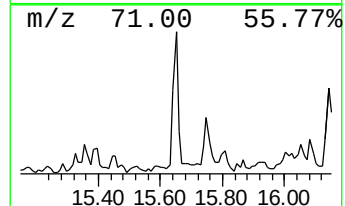
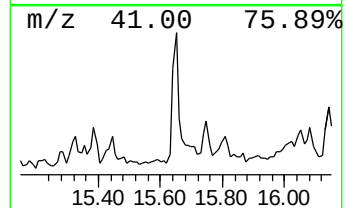
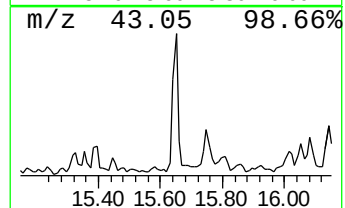
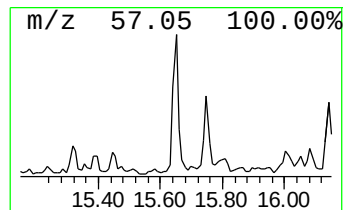
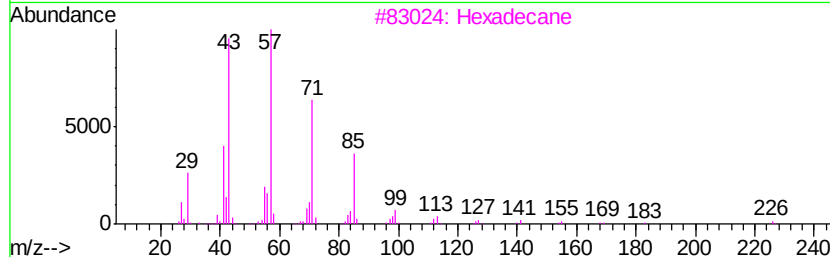
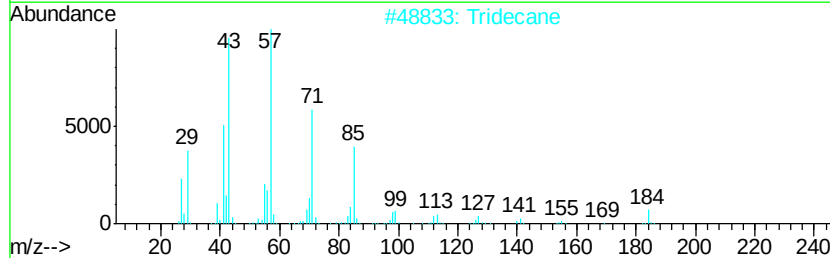
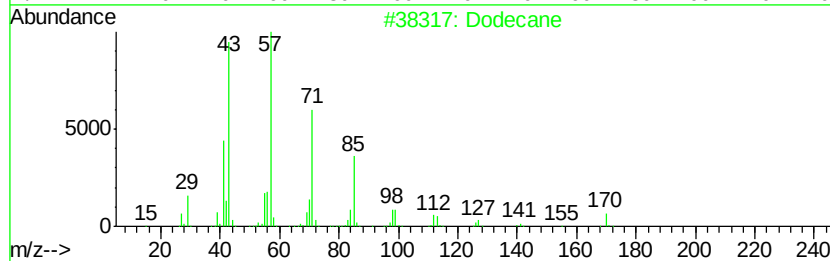
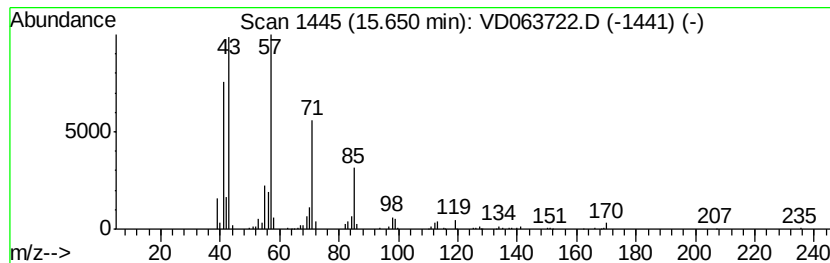
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	297.89 ug/l	1213850	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	95
2		Tridecane	184	C13H28	000629-50-5	90
3		Hexadecane	226	C16H34	000544-76-3	86
4		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	80
5		Undecane	156	C11H24	001120-21-4	72



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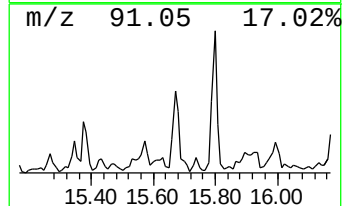
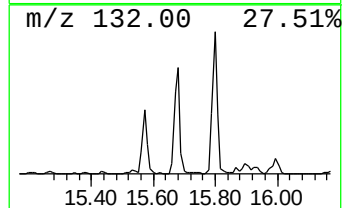
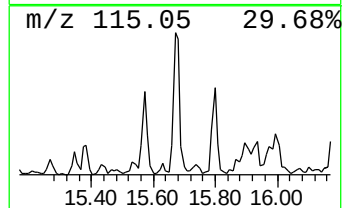
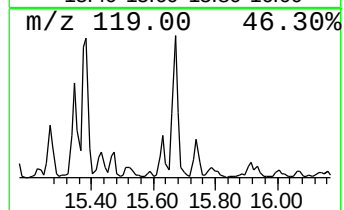
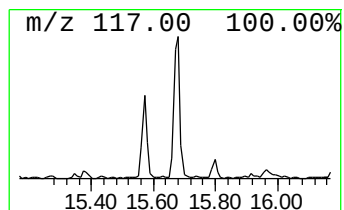
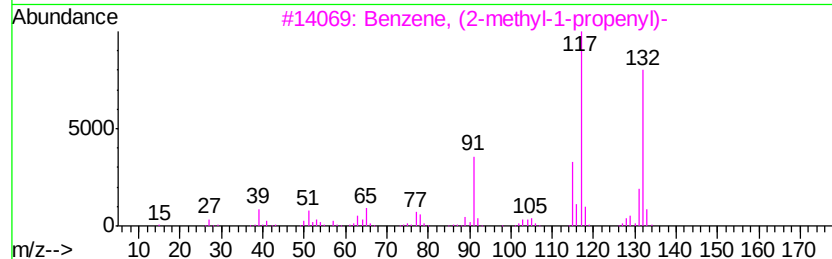
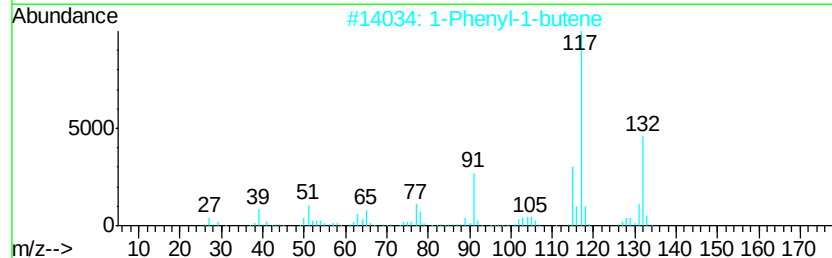
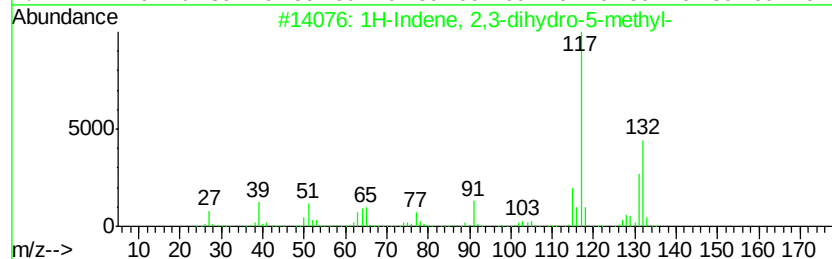
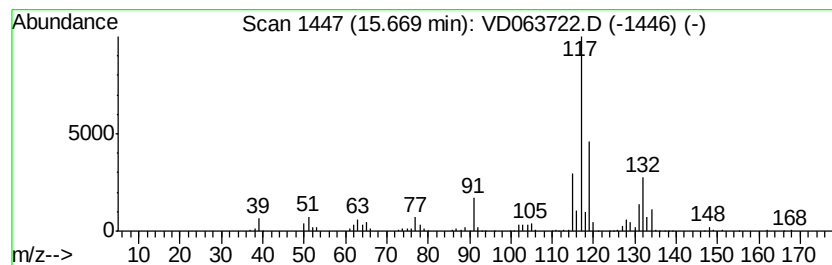
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	444.67 ug/l	1811950	1,4-Dichlorobenzene-d4	14.52

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	64
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD082119\
Data File : VD063722.D
Acq On : 21 Aug 2019 19:28
Operator : JC/SY
Sample : K4418-04
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 28 Sample Multiplier: 1

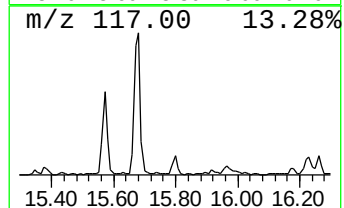
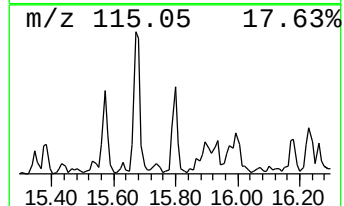
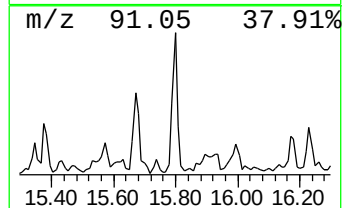
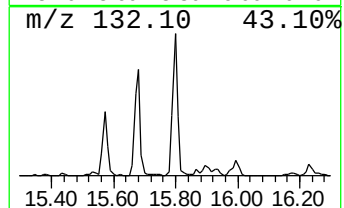
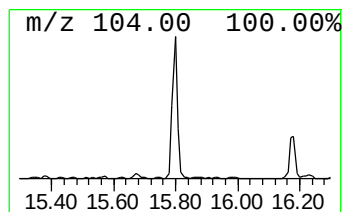
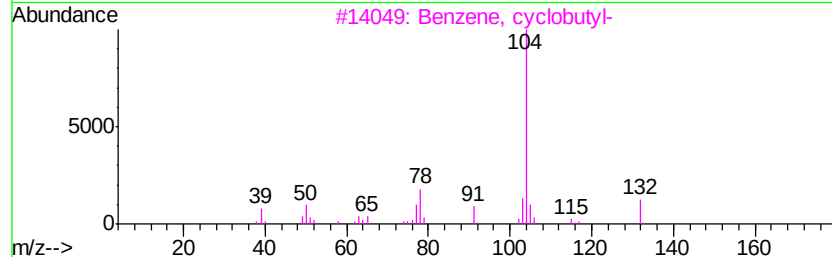
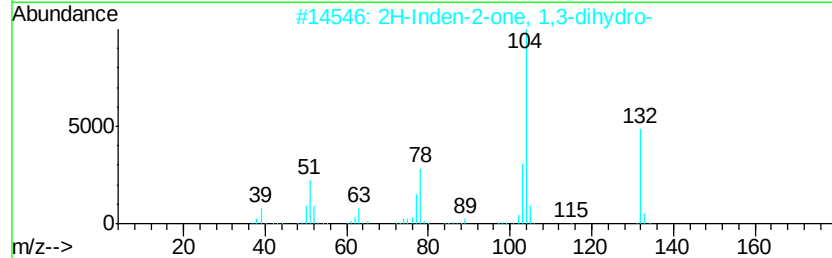
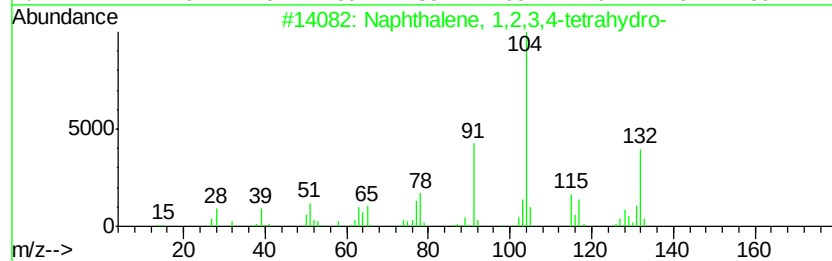
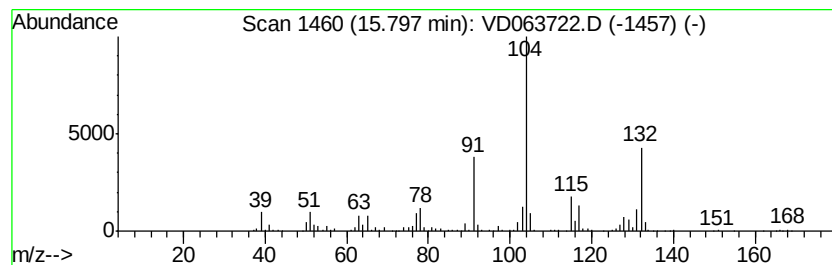
Instrument :
MSVOA_D
ClientSampled :
CFVP-228L

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	405.28 ug/l	1651470	1,4-Dichlorobenzene-d4	14.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 96
2		2H-Inden-2-one, 1,3-dihydro-	132 C9H8O	000615-13-4 53
3		Benzene, cyclobutyl-	132 C10H12	004392-30-7 53
4		Isoquinoline, 1,2,3,4-tetrahydro-	133 C9H11N	000091-21-4 43
5		Indan, epoxide	132 C9H8O	000768-22-9 41



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD082119\
Data File : VD063722.D
Acq On : 21 Aug 2019 19:28
Operator : JC/SY
Sample : K4418-04
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 28 Sample Multiplier: 1

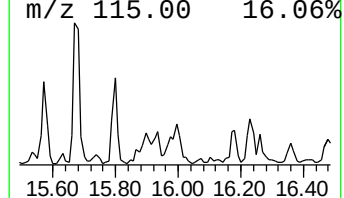
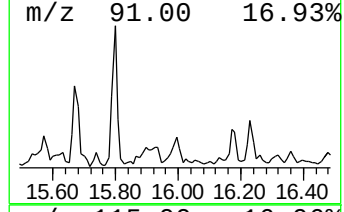
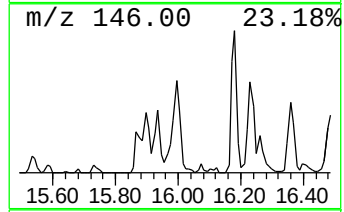
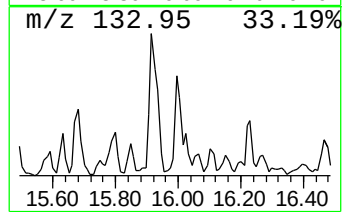
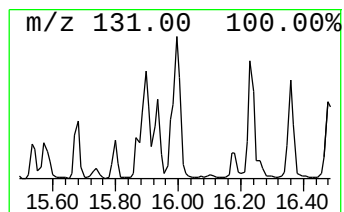
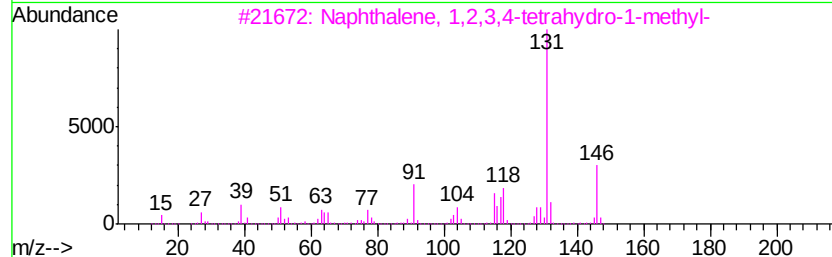
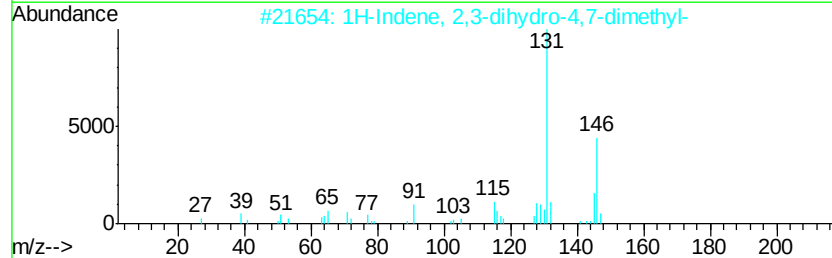
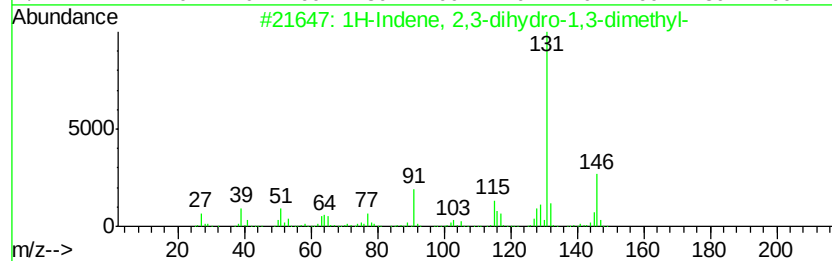
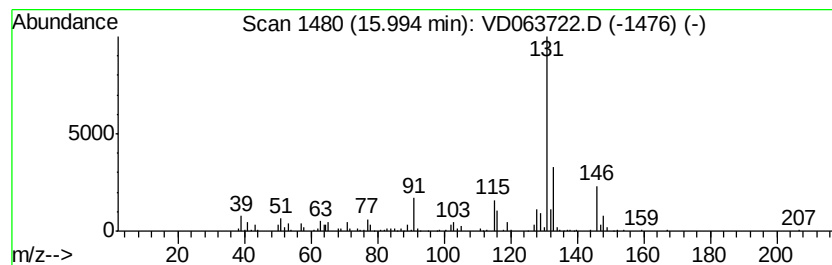
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	282.42 ug/l	1150800	1,4-Dichlorobenzene-d4	14.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	81
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	81
3	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	81
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	81
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD082119\
Data File : VD063722.D
Acq On : 21 Aug 2019 19:28
Operator : JC/SY
Sample : K4418-04
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 28 Sample Multiplier: 1

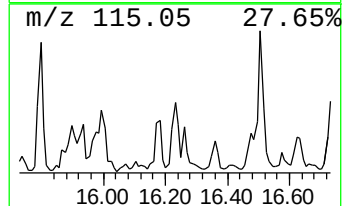
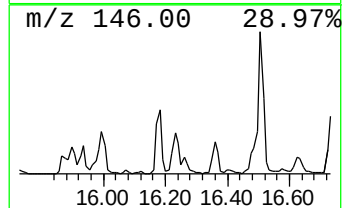
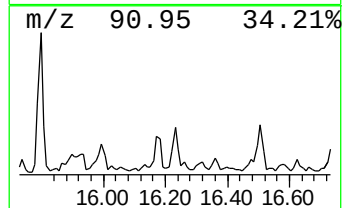
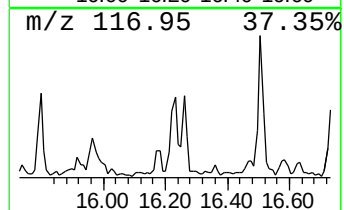
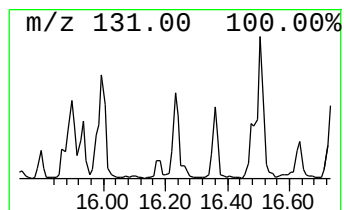
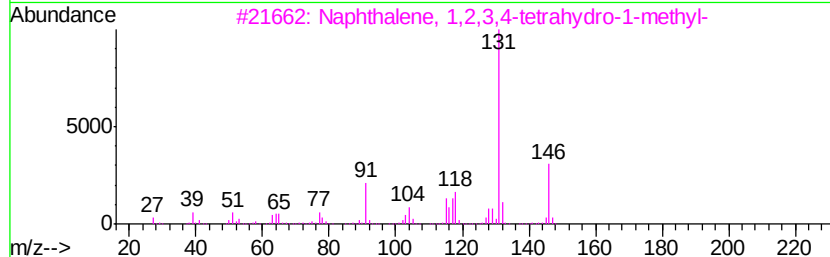
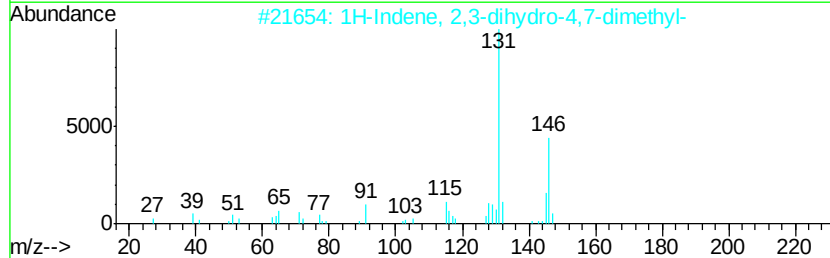
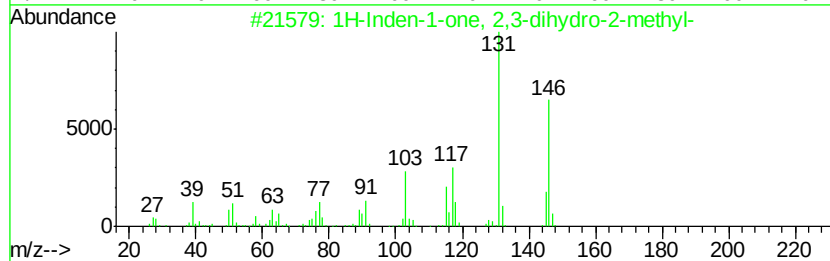
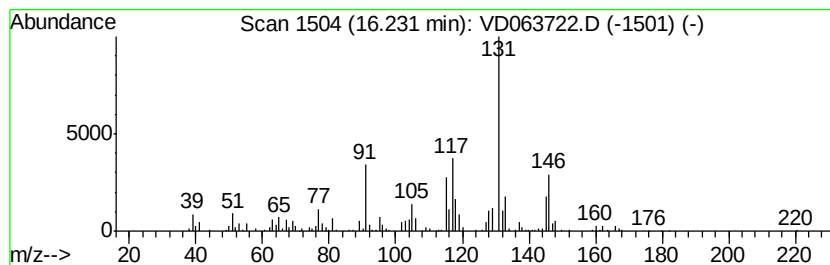
Instrument :
MSVOA_D
ClientSampleId :
CFVP-228L

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.23	354.45 ug/l	1444340	1,4-Dichlorobenzene-d4	14.52		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	87	
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81	
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	81	
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	76	
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD082119\
 Data File : VD063722.D
 Acq On : 21 Aug 2019 19:28
 Operator : JC/SY
 Sample : K4418-04
 Misc : 5.00µl/5ml/MSVOA D/SOIL
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CFVP-228L

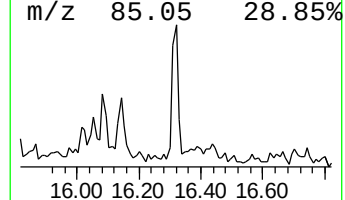
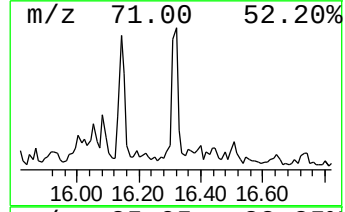
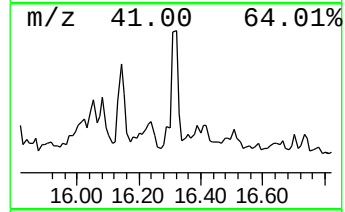
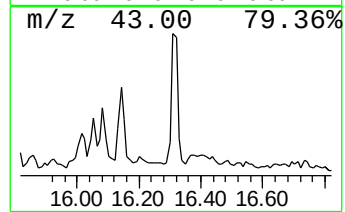
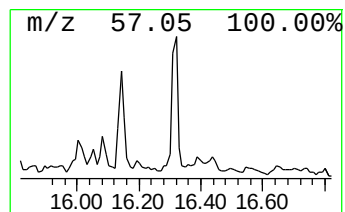
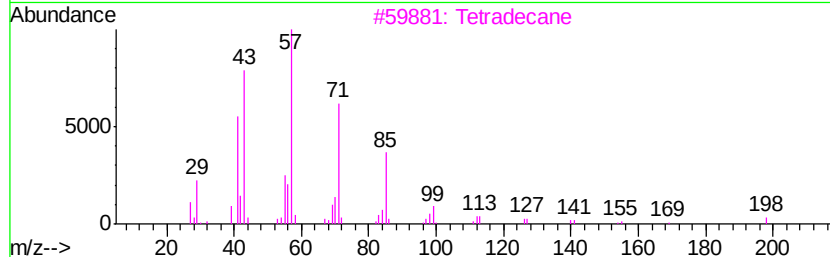
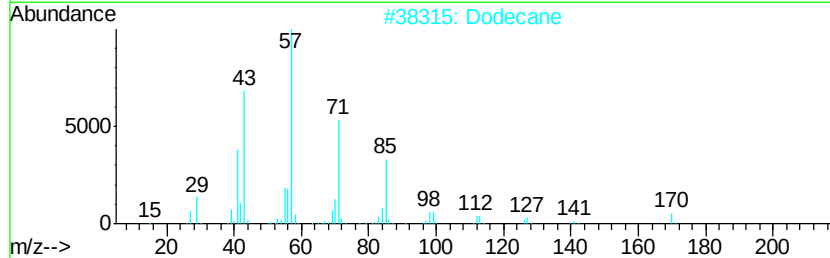
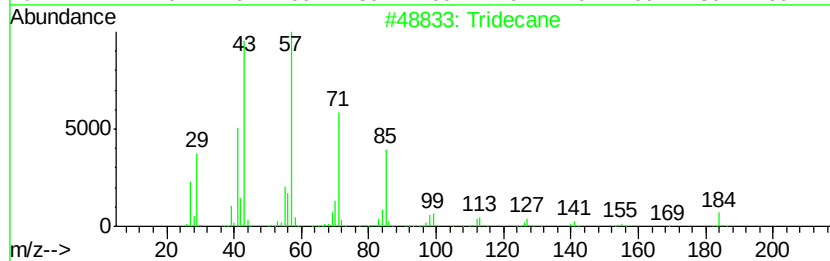
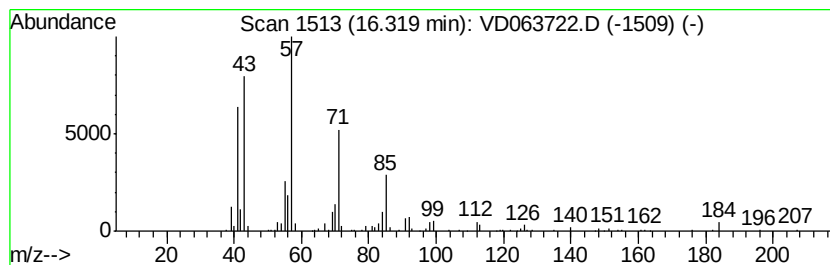
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 Quant Title : SW846 8260

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

 Peak Number 8 Tridecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.32	251.69 ug/l	1025610	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	87
2		Dodecane	170	C12H26	000112-40-3	86
3		Tetradecane	198	C14H30	000629-59-4	80
4		Undecane	156	C11H24	001120-21-4	80
5		Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD082119\
Data File : VD063722.D
Acq On : 21 Aug 2019 19:28
Operator : JC/SY
Sample : K4418-04
Misc : 5.00µl/5ml/MSVOA D/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CFVP-228L

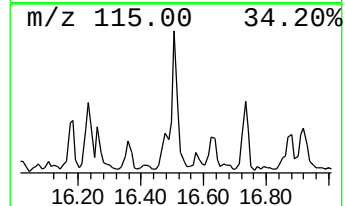
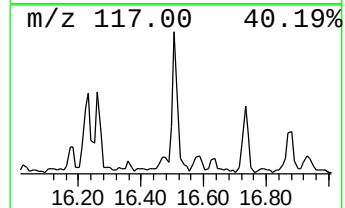
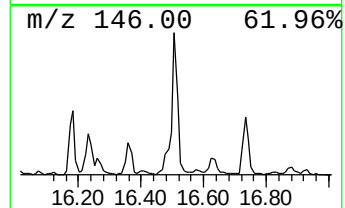
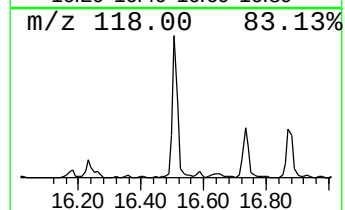
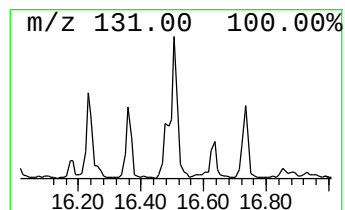
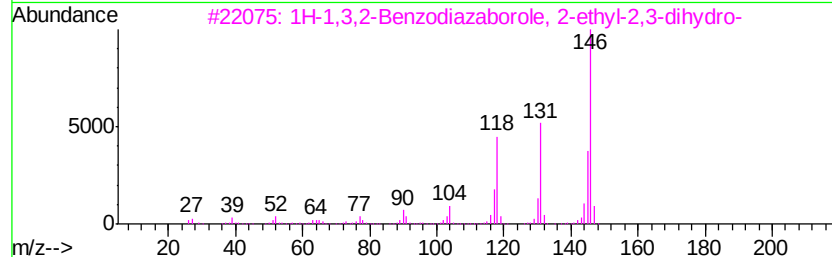
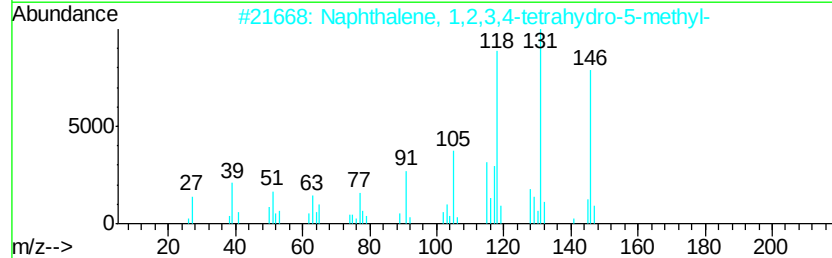
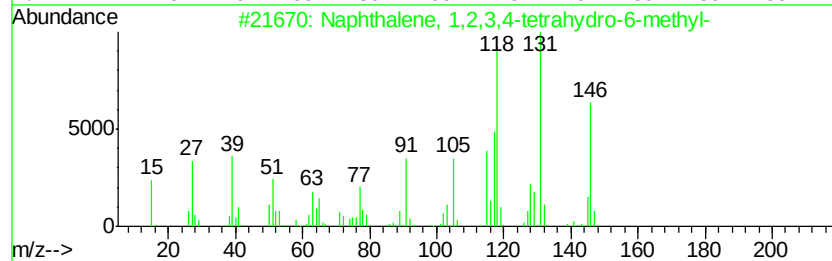
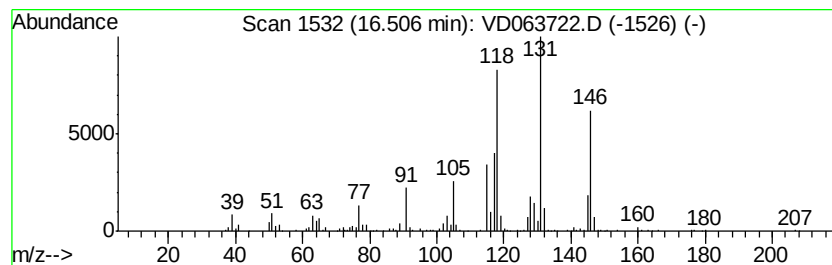
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	464.00 ug/l	1890740	1,4-Dichlorobenzene-d4	14.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	97
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3		1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	81
4		Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	74
5		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD082119\
Data File : VD063722.D
Acq On : 21 Aug 2019 19:28
Operator : JC/SY
Sample : K4418-04
Misc : 5.00g/5ml/MSVOA_D/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
CFVP-228L

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D080919S.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, 1-etheny...	15.14	232.5	ug/l	947520	4	14.52	203742	50.0
Benzene, 1-etheny...	15.57	307.7	ug/l	1253990	4	14.52	203742	50.0
Dodecane	15.65	297.9	ug/l	1213850	4	14.52	203742	50.0
1H-Indene, 2,3-di...	15.67	444.7	ug/l	1811950	4	14.52	203742	50.0
Naphthalene, 1,2,...	15.80	405.3	ug/l	1651470	4	14.52	203742	50.0
1H-Indene, 2,3-di...	15.99	282.4	ug/l	1150800	4	14.52	203742	50.0
1H-Inden-1-one, 2...	16.23	354.4	ug/l	1444340	4	14.52	203742	50.0
Tridecane	16.32	251.7	ug/l	1025610	4	14.52	203742	50.0
Naphthalene, 1,2,...	16.51	464.0	ug/l	1890740	4	14.52	203742	50.0