

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF052319\
 Data File : VF062725.D
 Acq On : 23 May 2019 14:29
 Operator : FY/SY
 Sample : K2976-04
 Misc : 6.22g/5mL/MSVOA F/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EP1-SB-E3-4R(17-20)

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_F\METHODS\82F051619S.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	4.320	377	384	398	rVB	467259	1515528	1.67%	0.086%
2	5.059	451	459	471	rBV2	943330	3194436	3.52%	0.182%
3	5.778	527	532	544	rVB	1164337	3070167	3.38%	0.175%
4	7.720	720	729	744	rBV	1268071	3626285	4.00%	0.206%
5	8.430	794	801	808	rBV	749116	2175321	2.40%	0.124%
6	8.607	808	819	827	rBV2	1271271	4667088	5.14%	0.265%
7	8.775	831	836	849	rBV	1084349	4186267	4.61%	0.238%
8	9.120	864	871	876	rBV	2845459	9079333	10.01%	0.516%
9	9.248	880	884	892	rVB	2322876	6253164	6.89%	0.356%
10	9.426	892	902	908	rBV	2807972	8631105	9.51%	0.491%
11	9.554	908	915	920	rBV4	428733	1362558	1.50%	0.077%
12	9.682	920	928	935	rBV5	1253142	5118092	5.64%	0.291%
13	9.840	935	944	947	rBV2	1947631	7567111	8.34%	0.430%
14	9.918	947	952	955	rBV2	7353372	20135464	22.20%	1.145%
15	10.125	969	973	977	rBV2	1045053	2599680	2.87%	0.148%
16	10.214	977	982	985	rBV	1663714	5069485	5.59%	0.288%
17	10.263	985	987	992	rVB	2338568	4644309	5.12%	0.264%
18	10.382	992	999	1005	rBV2	9156467	39407590	43.44%	2.240%
19	10.490	1005	1010	1015	rVB2	7651567	20699208	22.82%	1.177%
20	10.648	1015	1026	1034	rBV6	12325888	65951262	72.70%	3.749%
21	10.806	1036	1042	1049	rVV4	24061463	90714220	100.00%	5.157%
22	10.914	1049	1053	1054	rVV2	8851368	21535909	23.74%	1.224%
23	10.963	1054	1058	1061	rVV	12893778	37050422	40.84%	2.106%
24	11.072	1065	1069	1072	rVB2	7010624	16200172	17.86%	0.921%
25	11.170	1072	1079	1082	rBV4	16207246	55676469	61.38%	3.165%
26	11.239	1082	1086	1091	rVB3	15845680	39863561	43.94%	2.266%
27	11.387	1091	1101	1105	rVB5	18631634	68061237	75.03%	3.869%
28	11.446	1105	1107	1109	rBV3	1269032	1785151	1.97%	0.101%
29	11.496	1109	1112	1114	rBV2	7315693	15474562	17.06%	0.880%
30	11.555	1114	1118	1120	rVV3	11703940	32465530	35.79%	1.846%
31	11.594	1120	1122	1125	rVV4	11407425	24287230	26.77%	1.381%
32	11.663	1125	1129	1132	rVV3	20354722	51858417	57.17%	2.948%
33	11.722	1132	1135	1138	rVV2	14559912	32444390	35.77%	1.845%
34	11.772	1138	1140	1142	rVV2	7894174	14912531	16.44%	0.848%

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35	11.821	1142	1145	1148	rVV2	10138079	29368618	32.37%	1.670%
36	11.910	1152	1154	1158	rVB2	12662625	23461999	25.86%	1.334%
37	11.979	1158	1161	1163	rBV2	10137941	20179158	22.24%	1.147%
38	12.048	1163	1168	1169	rBV2	11352442	26888020	29.64%	1.529%
39	12.097	1169	1173	1177	rVB4	24743654	61543186	67.84%	3.499%
40	12.166	1177	1180	1184	rBV3	10831980	23566493	25.98%	1.340%
41	12.245	1184	1188	1189	rBV2	11898324	24500315	27.01%	1.393%
42	12.294	1189	1193	1197	rVB2	35944975	88003880	97.01%	5.003%
43	12.373	1197	1201	1204	rBV3	8699279	19459823	21.45%	1.106%
44	12.452	1207	1209	1211	rBV2	5450096	7531617	8.30%	0.428%
45	12.570	1219	1221	1223	rVB2	6090834	7612972	8.39%	0.433%
46	12.688	1229	1233	1237	rBV2	26775208	58760015	64.77%	3.341%
47	12.757	1237	1240	1241	rVV	6857387	13003852	14.33%	0.739%
48	12.807	1241	1245	1247	rVV2	20089469	45817821	50.51%	2.605%
49	12.866	1247	1251	1253	rVV2	14938249	38436254	42.37%	2.185%
50	12.905	1253	1255	1259	rVV3	11040445	22680215	25.00%	1.289%
51	12.994	1259	1264	1269	rVV3	13930303	37144040	40.95%	2.112%
52	13.073	1269	1272	1275	rVB2	9557222	20450189	22.54%	1.163%
53	13.161	1275	1281	1284	rBV3	21100768	70377966	77.58%	4.001%
54	13.221	1284	1287	1292	rVV3	26940768	70850261	78.10%	4.028%
55	13.299	1292	1295	1297	rVV2	15191866	30081522	33.16%	1.710%
56	13.349	1297	1300	1304	rVV2	17163861	40404217	44.54%	2.297%
57	13.437	1307	1309	1312	rVV	14372835	20187335	22.25%	1.148%
58	13.507	1314	1316	1318	rVV2	4314648	7979887	8.80%	0.454%
59	13.556	1318	1321	1323	rVV	16572232	29717143	32.76%	1.689%
60	13.595	1323	1325	1332	rVB4	24260011	49878170	54.98%	2.836%
61	13.684	1332	1334	1338	rBV	7257176	11660447	12.85%	0.663%
62	13.753	1338	1341	1344	rBV	11386126	17517901	19.31%	0.996%
63	13.802	1344	1346	1349	rVB	4057604	5730428	6.32%	0.326%
64	13.861	1349	1352	1353	rBV	6855582	10340262	11.40%	0.588%
65	13.921	1353	1358	1361	rVV3	19463001	55505619	61.19%	3.156%
66	13.999	1364	1366	1370	rVB3	1636163	2828883	3.12%	0.161%
67	14.098	1373	1376	1377	rBV2	2260179	4129056	4.55%	0.235%
68	14.118	1377	1378	1381	rVB	2913769	3191659	3.52%	0.181%
69	14.197	1381	1386	1389	rBV3	6610041	15767978	17.38%	0.896%
70	14.256	1390	1392	1398	rVB	4537218	6834646	7.53%	0.389%
71	14.335	1398	1400	1403	rVB2	2198001	2870528	3.16%	0.163%

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72	14.394	1403	1406	1409	rVB2	1673257	2880114	3.17%	0.164%
73	14.453	1409	1412	1415	rBV3	500847	1144189	1.26%	0.065%
74	14.522	1415	1419	1424	rVB4	1085740	3575553	3.94%	0.203%
75	14.768	1441	1444	1450	rVB3	1409451	2251849	2.48%	0.128%
76	15.025	1467	1470	1475	rVB2	605782	1037382	1.14%	0.059%
77	15.310	1493	1499	1503	rBV2	838344	1408710	1.55%	0.080%
78	15.439	1508	1512	1521	rVB	698370	1106756	1.22%	0.063%

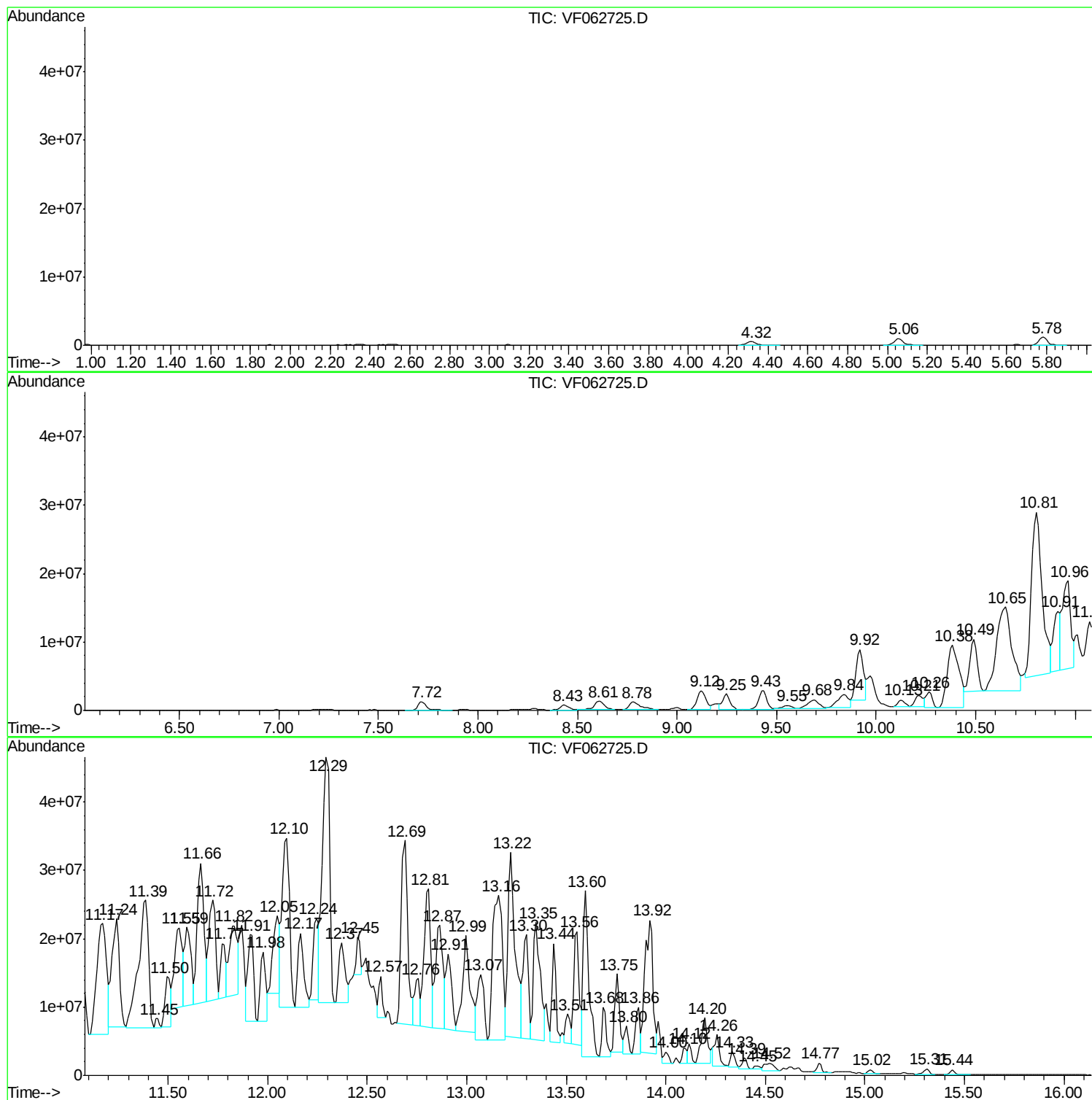
Sum of corrected areas: 1758950182

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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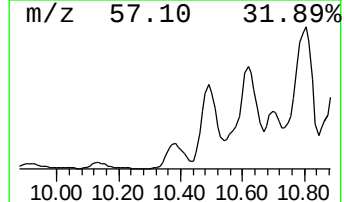
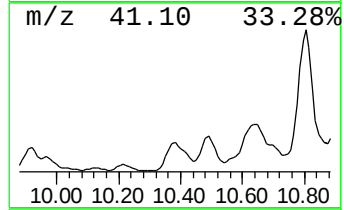
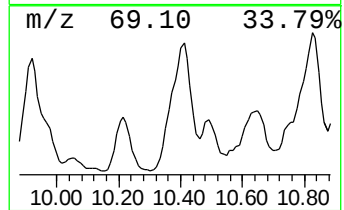
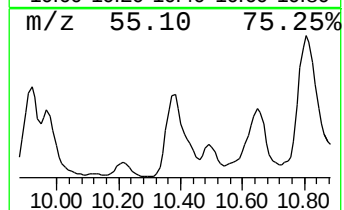
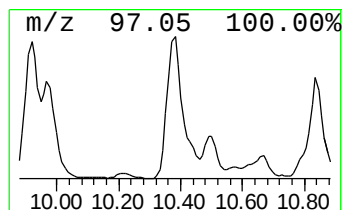
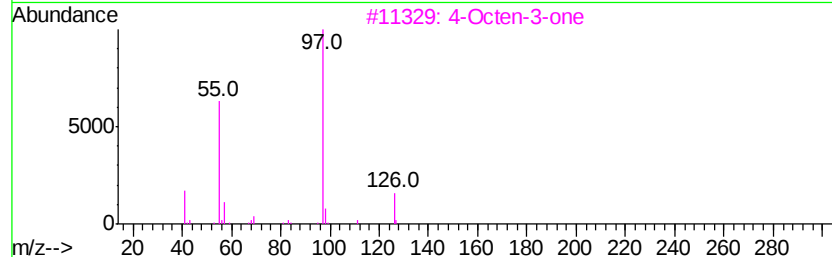
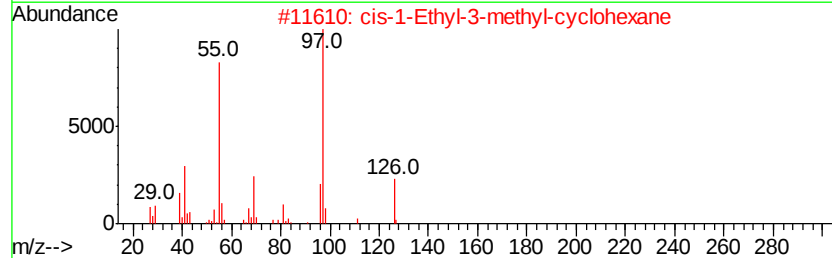
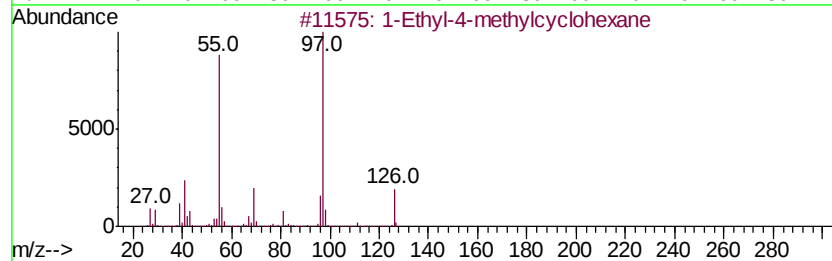
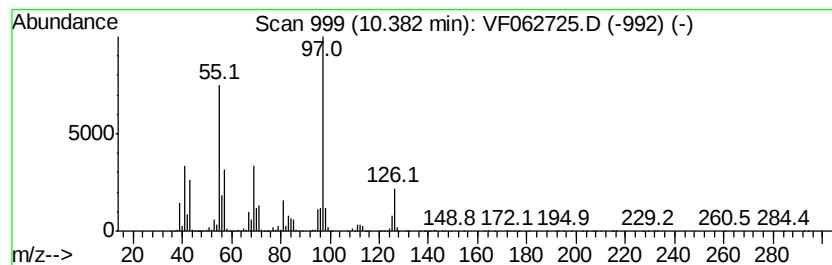
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Ethyl-4-methylcyclohexane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	97.86 ug/l	39407600	Chlorobenzene-d5	9.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	83
2		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	72
3		4-Octen-3-one	126	C8H14O	014129-48-7	72
4		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	72
5		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	70



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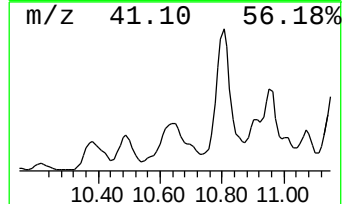
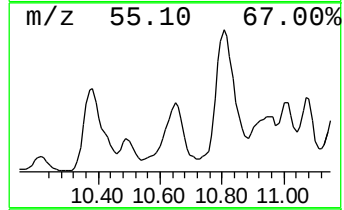
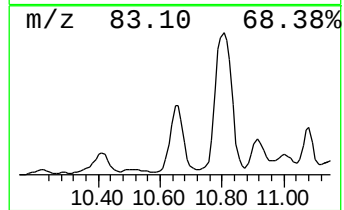
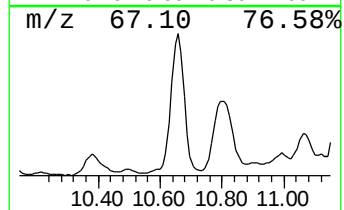
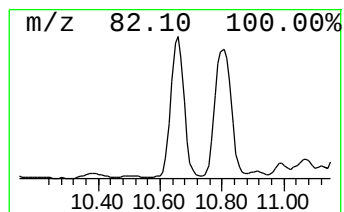
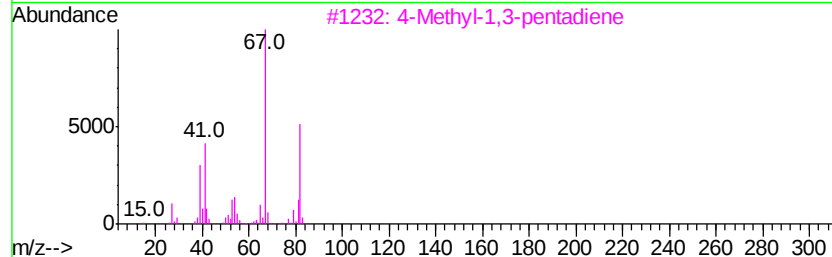
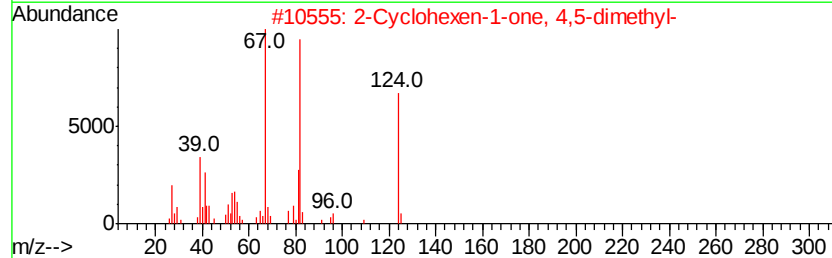
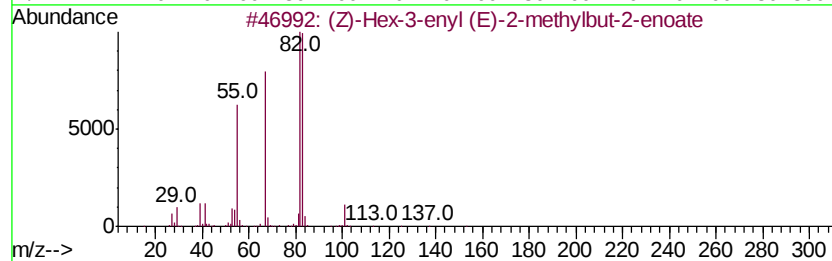
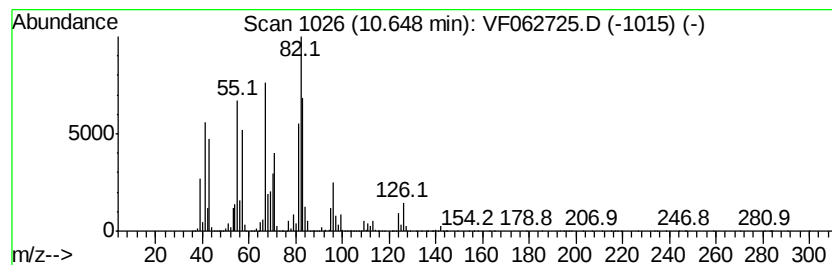
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown10.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	163.77 ug/l	65951300	Chlorobenzene-d5	9.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(Z)-Hex-3-enyl (E)-2-methylbut-2...	182	C11H18O2	1000373-73-0	47
2		2-Cyclohexen-1-one, 4,5-dimethyl-	124	C8H12O	005715-25-3	38
3		4-Methyl-1,3-pentadiene	82	C6H10	000926-56-7	38
4		3-Aminocrotononitrile	82	C4H6N2	001118-61-2	38
5		Bicyclo[2.2.1]heptane, 2-ethyl-	124	C9H16	002146-41-0	30



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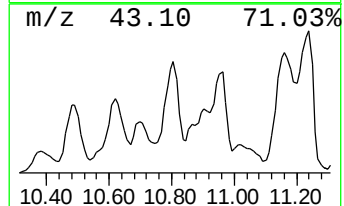
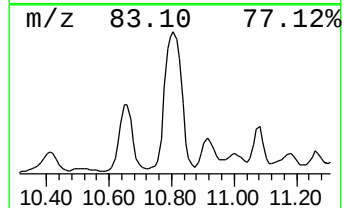
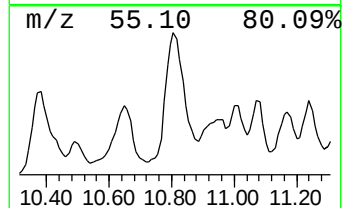
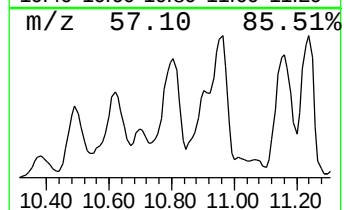
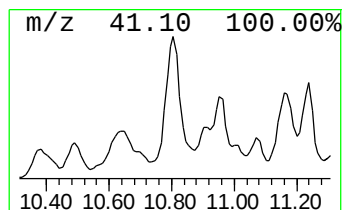
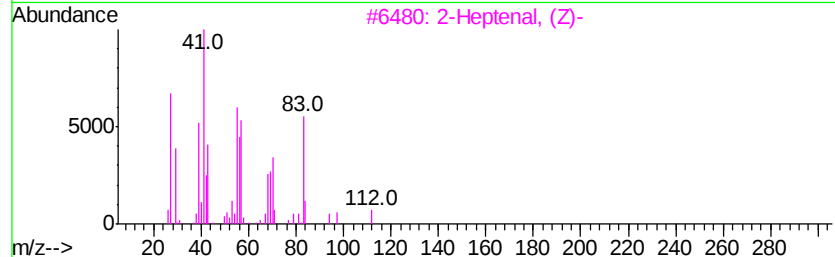
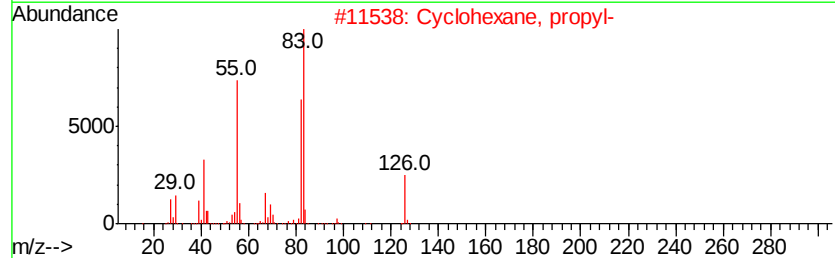
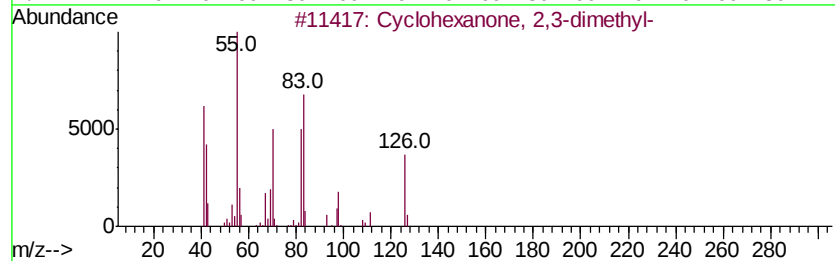
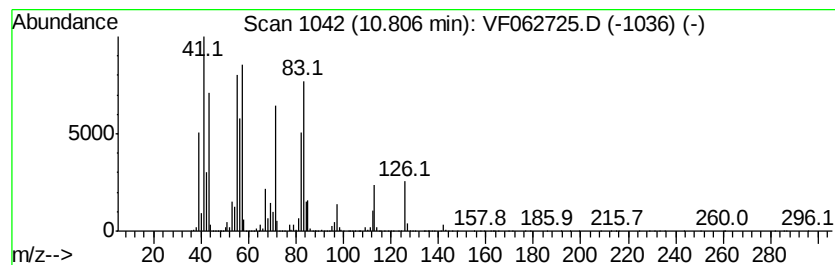
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown10.81 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.81	225.26 ug/l	90714200	Chlorobenzene-d5	9.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	46
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	46
3	2-Heptenal, (Z)-	112	C7H12O	057266-86-1	45
4	2-Heptenal, (E)-	112	C7H12O	018829-55-5	45
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	43



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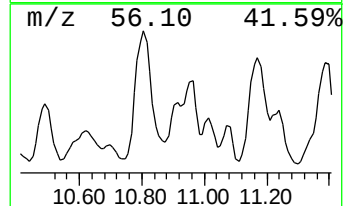
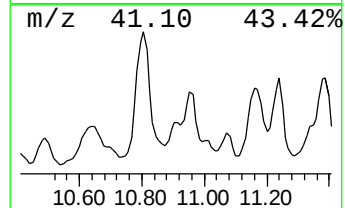
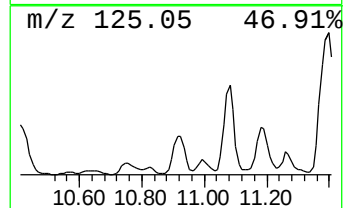
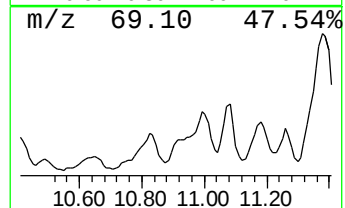
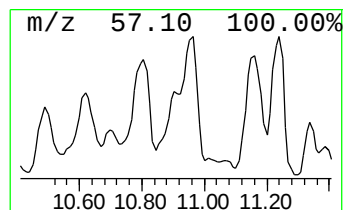
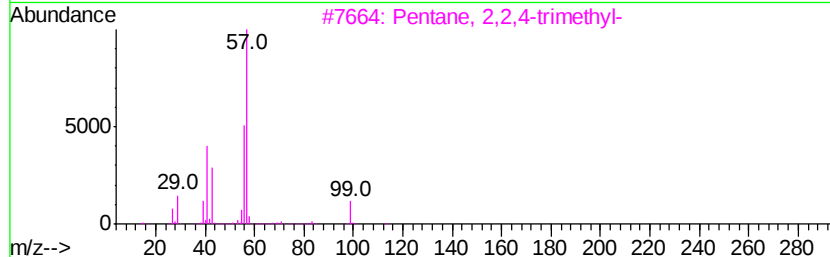
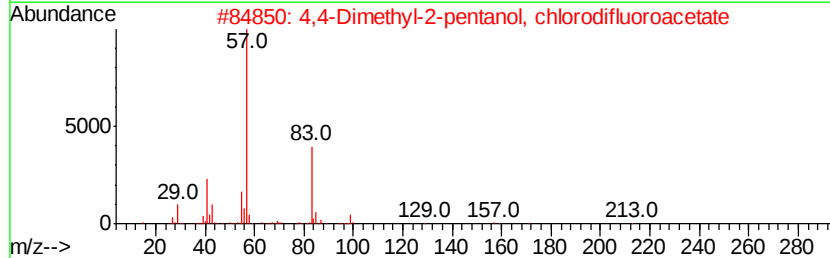
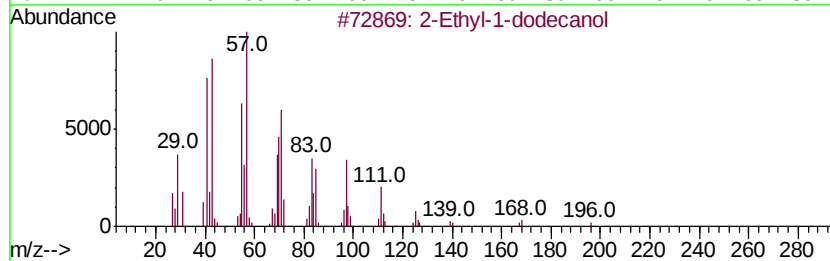
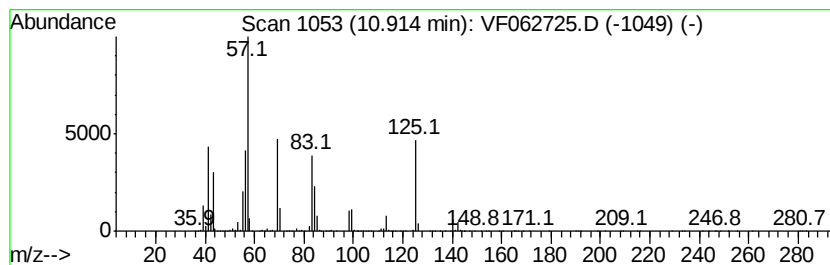
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Quant Title : SW846 8260

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

Peak Number 4 unknown10.91 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	53.48 ug/l	21535900	Chlorobenzene-d5	9.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Ethyl-1-dodecanol	214	C14H30O	019780-33-7	45
2		4,4-Dimethyl-2-pentanol, chlorod...	228	C9H15ClF2O2	1000376-27-0	27
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	22
4		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	14
5		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	14



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
Operator : FY/SY
Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

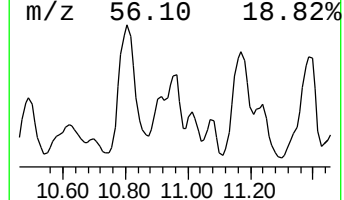
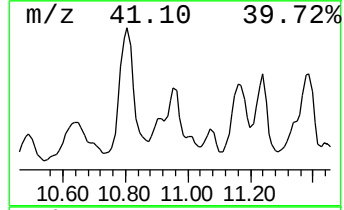
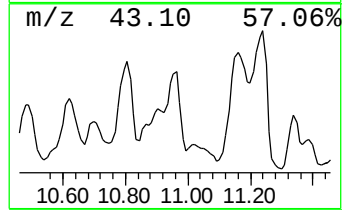
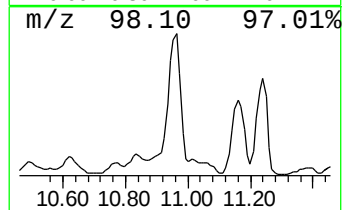
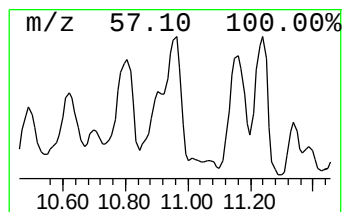
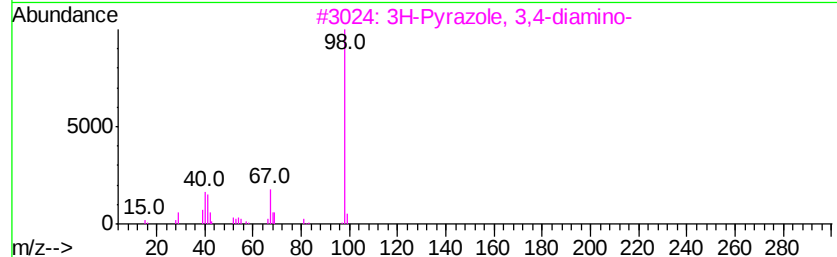
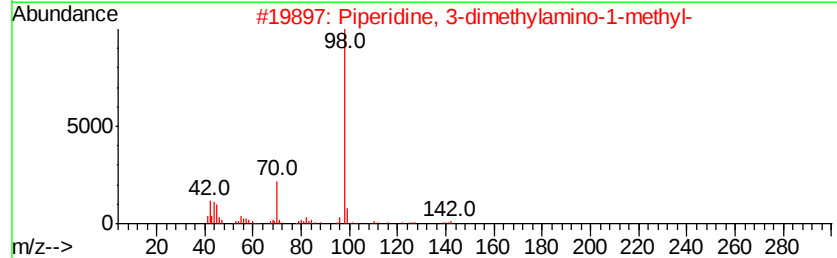
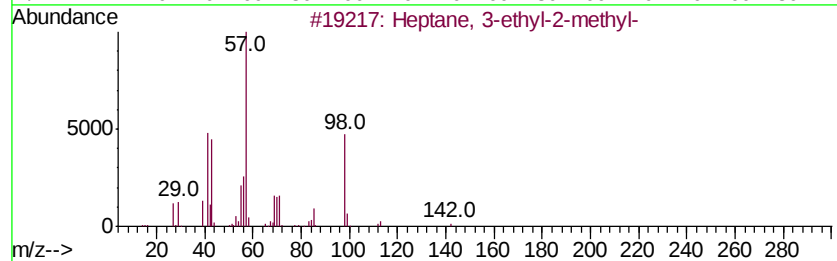
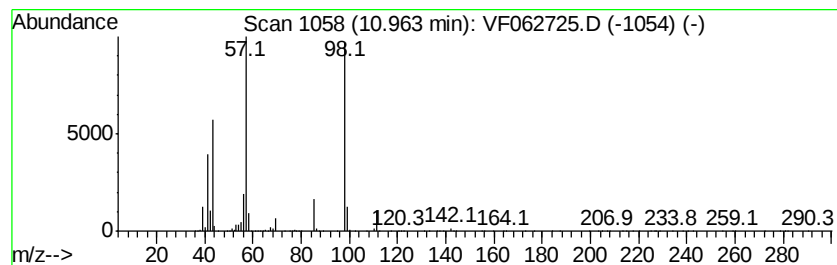
Instrument :
MSVOA_F
ClientSampleId :
EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 5 Heptane, 3-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.96	92.00 ug/l	37050400	Chlorobenzene-d5	9.92		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
2		Piperidine, 3-dimethylamino-1-me...	142	C8H18N2	1000126-80-8	47
3		3H-Pyrazole, 3,4-diamino-	98	C3H6N4	1000288-40-0	47
4		Heptane, 4-propyl-	142	C10H22	003178-29-8	43
5		Benzamide, 2-hydroxy-4-propoxy-N...	306	C17H26N2O3	1000295-50-1	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
Operator : FY/SY
Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

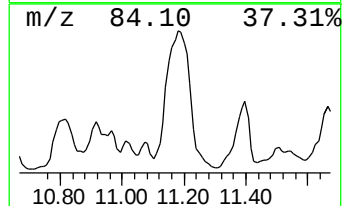
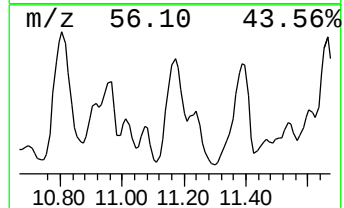
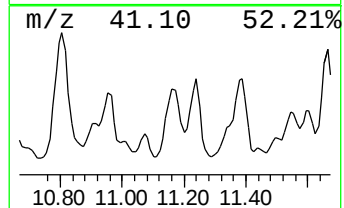
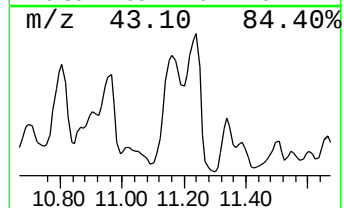
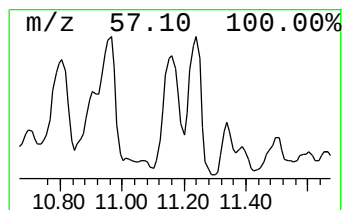
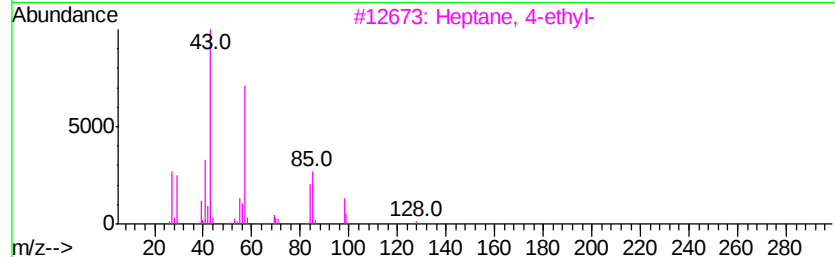
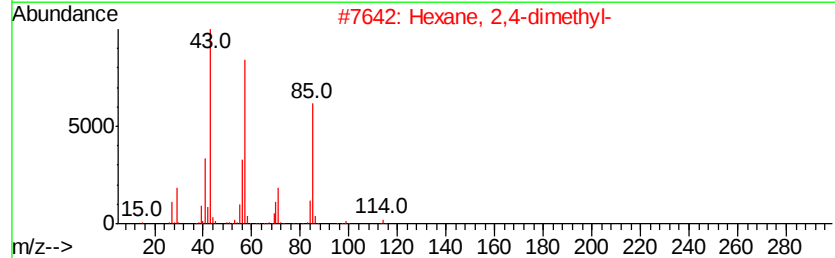
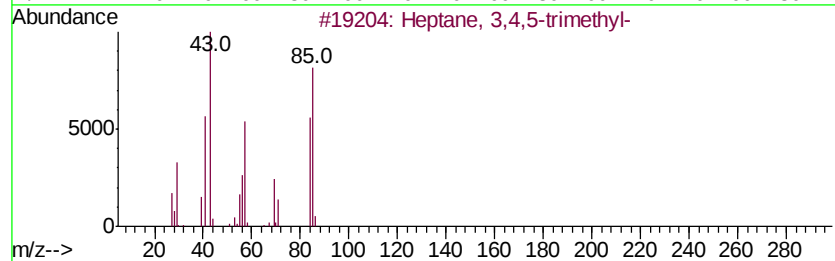
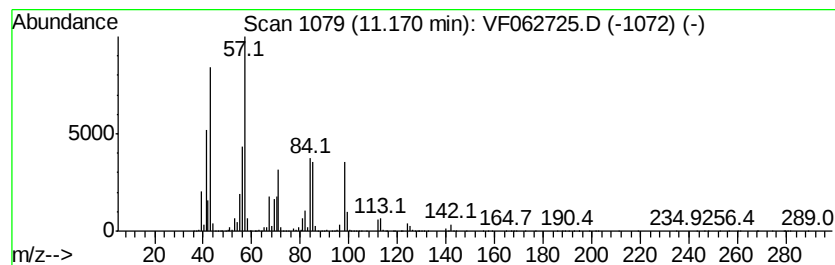
Instrument :
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ClientSampleId :
EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 6 Heptane, 3,4,5-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.17	138.26 ug/l	55676500	Chlorobenzene-d5	9.92		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	52	
2	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	47	
3	Heptane, 4-ethyl-	128	C9H20	002216-32-2	43	
4	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	43	
5	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	38	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
Operator : FY/SY
Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

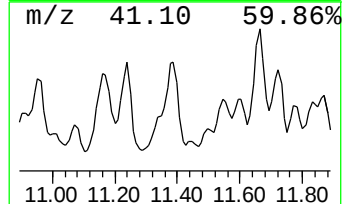
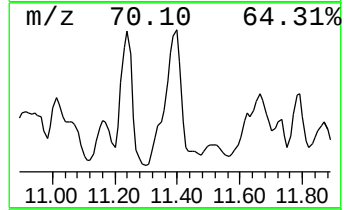
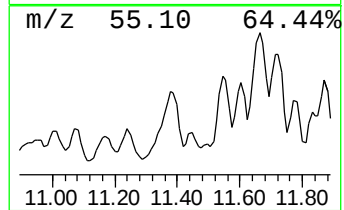
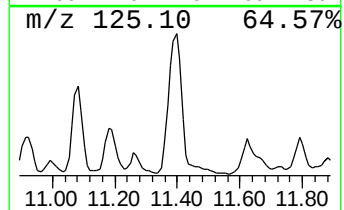
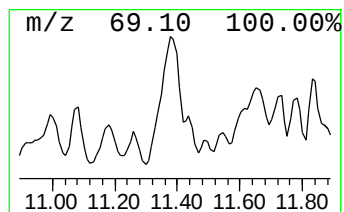
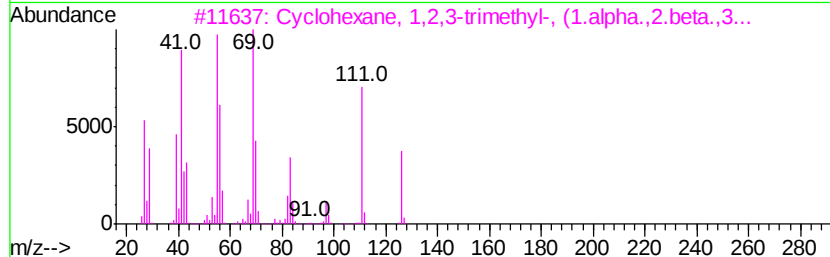
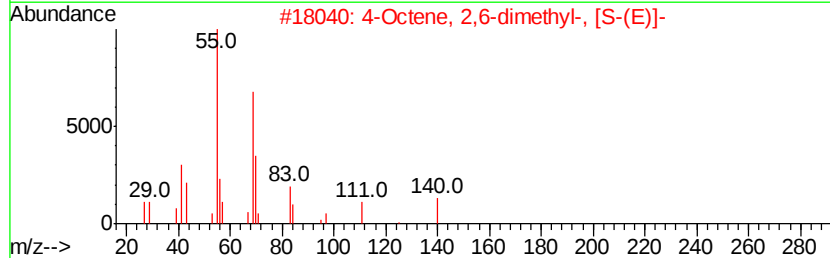
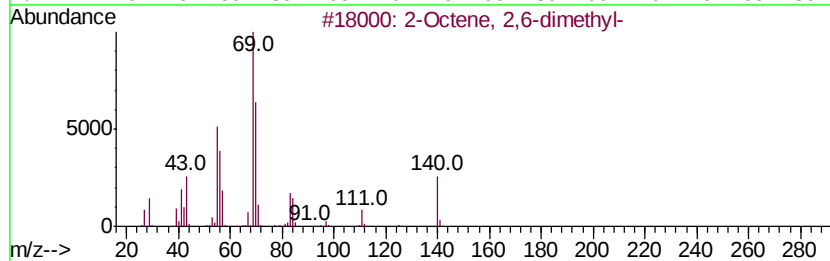
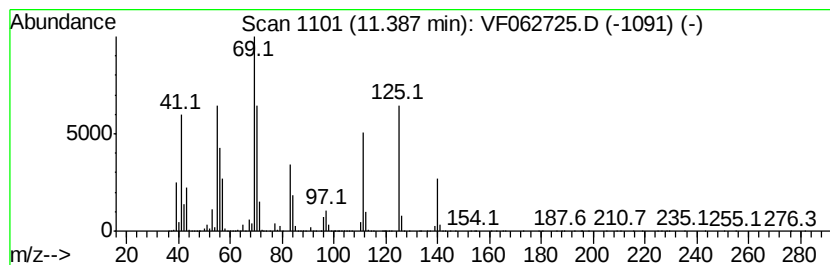
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ClientSampleId :
EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 7 2-Octene, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.39	57.91 ug/l	68061200	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Octene, 2,6-dimethyl-	140 C10H20	004057-42-5	91
2	4-Octene, 2,6-dimethyl-, [S-(E)]-	140 C10H20	062960-76-3	68
3	Cyclohexane, 1,2,3-trimethyl-, (...)	126 C9H18	001678-81-5	62
4	4-Octene, 2,6-dimethyl-, [S-(Z)]-	140 C10H20	062960-77-4	58
5	cis-1-Butyl-2-methylcyclopropane	112 C8H16	038851-69-3	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
Operator : FY/SY
Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EP1-SB-E3-4R(17-20)

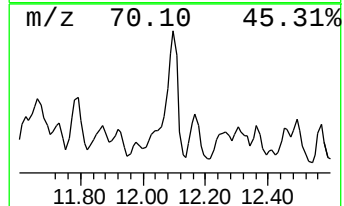
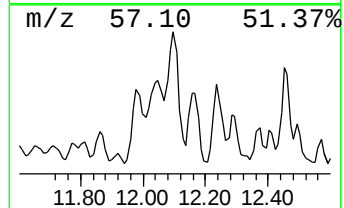
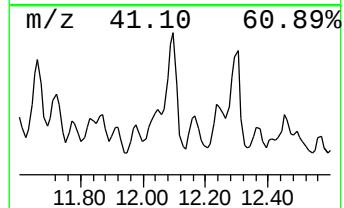
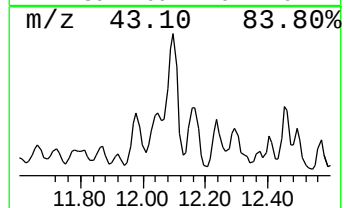
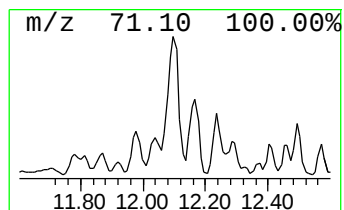
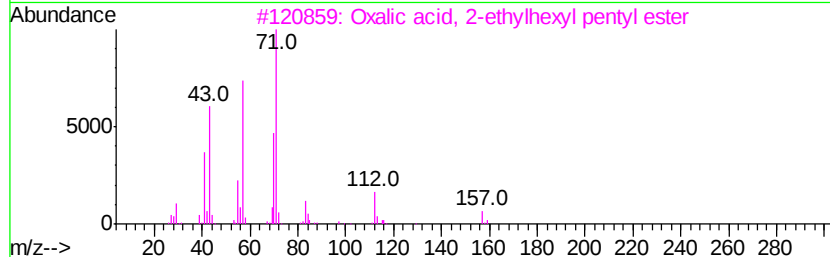
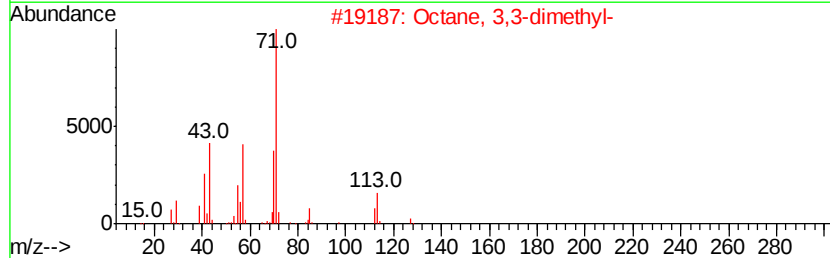
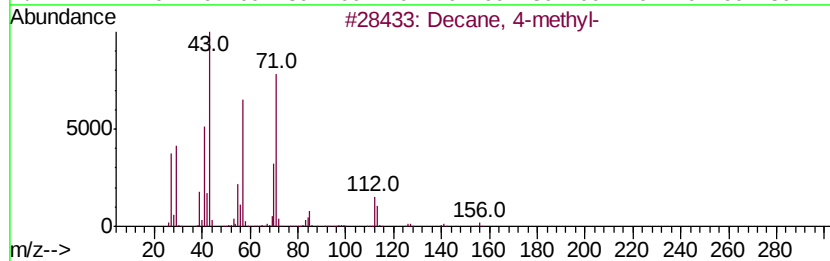
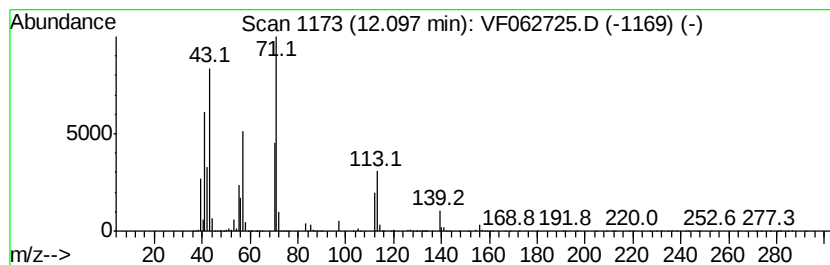
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 8 Decane, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	52.37 ug/l	61543200	1,4-Dichlorobenzene-d4	12.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 4-methyl-	156	C11H24	002847-72-5	72
2		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	59
3		Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	59
4		Oxalic acid, butyl 2-ethylhexyl ...	258	C14H26O4	1000309-38-6	59
5		Sulfurous acid, octyl 2-pentyl e...	264	C13H28O3S	1000309-15-7	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
Operator : FY/SY
Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

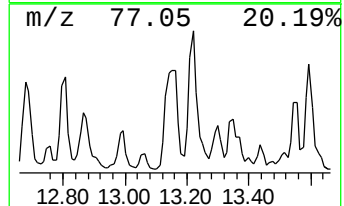
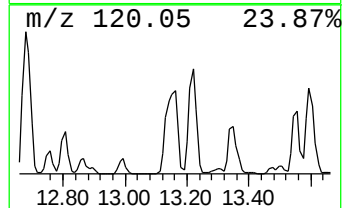
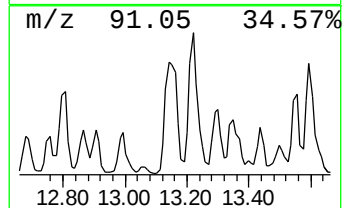
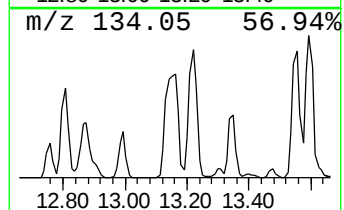
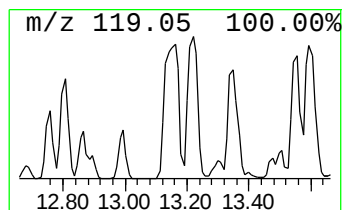
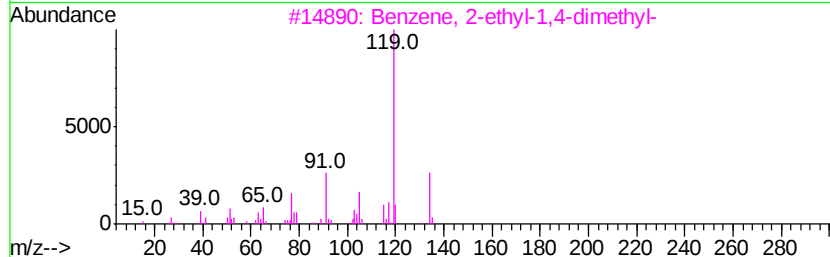
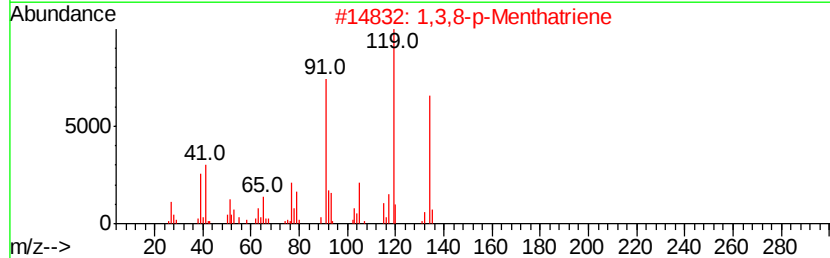
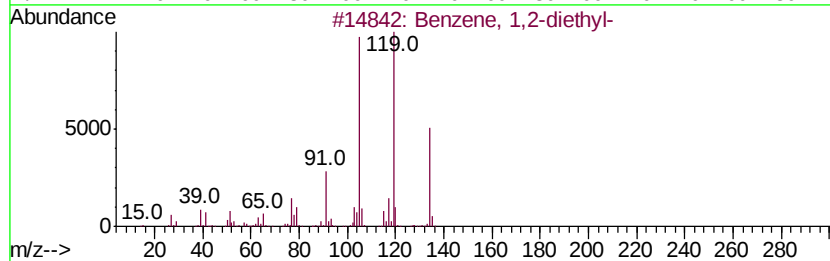
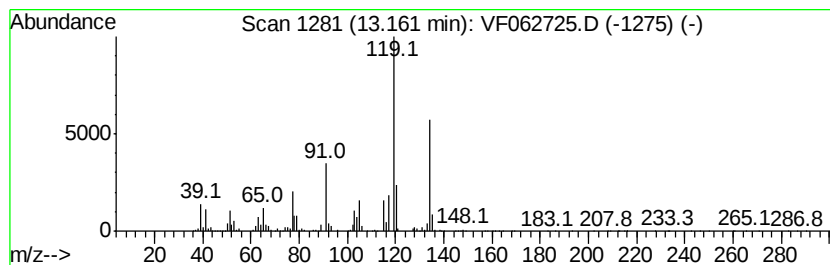
Instrument :
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ClientSampleId :
EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	59.89 ug/l	70378000	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2-diethyl-	134 C10H14	000135-01-3 90
2		1,3,8-p-Menthatriene	134 C10H14	018368-95-1 83
3		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 81
4		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 74
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134 C10H14	076089-59-3 68



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF052319\
Data File : VF062725.D
Acq On : 23 May 2019 14:29
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Sample : K2976-04
Misc : 6.22g/5mL/MSVOA F/SOIL
ALS Vial : 9 Sample Multiplier: 1

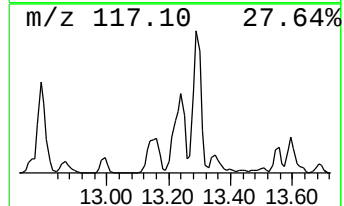
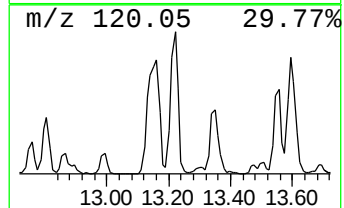
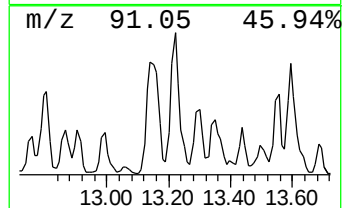
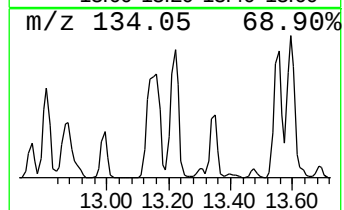
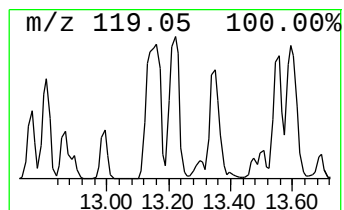
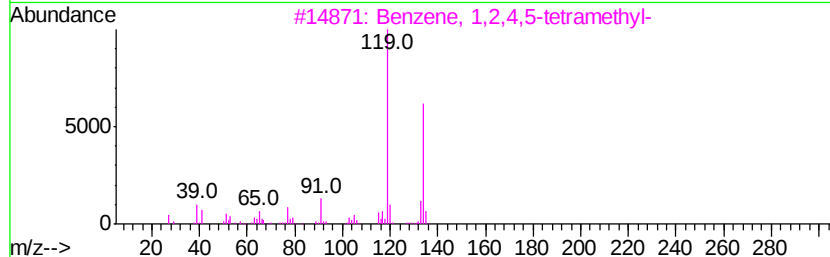
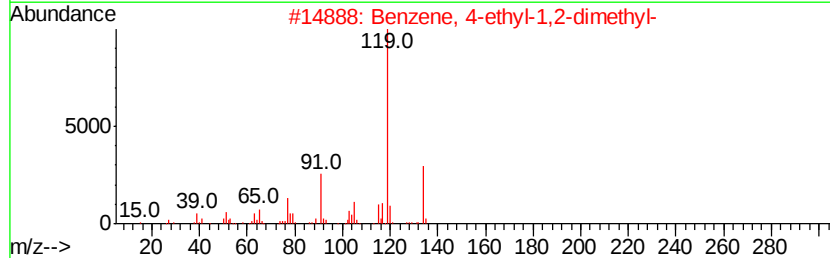
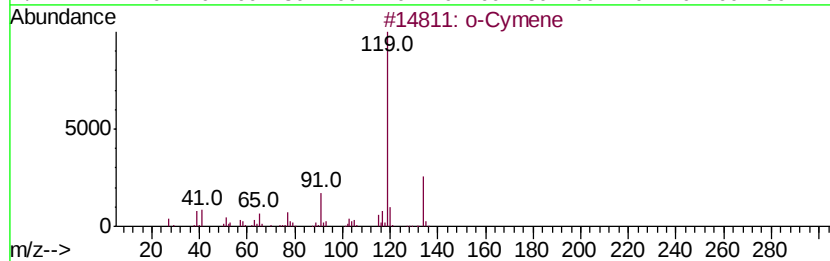
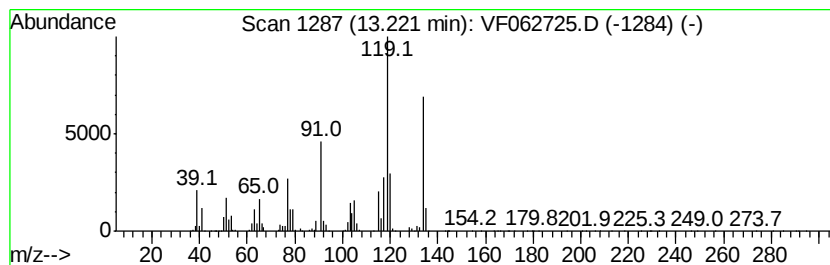
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MSVOA_F
ClientSampleId :
EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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

Peak Number 10 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	60.29 ug/l	70850300	1,4-Dichlorobenzene-d4	12.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 64
2		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 60
3		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 58
4		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 58
5		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 58



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_F\DATA\VF052319\
 Data File : VF062725.D
 Acq On : 23 May 2019 14:29
 Operator : FY/SY
 Sample : K2976-04
 Misc : 6.22g/5mL/MSV0A_F/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSV0A_F
 ClientSampleId :
 EP1-SB-E3-4R(17-20)

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_F\METHODS\82F051619S.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
1-Ethyl-4-methylc...	10.38	97.9	ug/l	39407600	3	9.92	20135500	50.0
unknown10.65	10.65	163.8	ug/l	65951300	3	9.92	20135500	50.0
unknown10.81	10.81	225.3	ug/l	90714200	3	9.92	20135500	50.0
unknown10.91	10.91	53.5	ug/l	21535900	3	9.92	20135500	50.0
Heptane, 3-ethyl-...	10.96	92.0	ug/l	37050400	3	9.92	20135500	50.0
Heptane, 3,4,5-tr...	11.17	138.3	ug/l	55676500	3	9.92	20135500	50.0
2-Octene, 2,6-dim...	11.39	57.9	ug/l	68061200	4	12.63	58760000	50.0
Decane, 4-methyl-	12.10	52.4	ug/l	61543200	4	12.63	58760000	50.0
Benzene, 1,2-diet...	13.16	59.9	ug/l	70378000	4	12.63	58760000	50.0
o-Cymene	13.22	60.3	ug/l	70850300	4	12.63	58760000	50.0