

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY061920\
 Data File : VY002958.D
 Acq On : 19 Jun 2020 23:22
 Operator : SY/MD
 Sample : L3063-15
 Misc : 5.06G/5ML/MSVOA Y/SOIL
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TW-11

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_Y\METHODS\82Y061920S.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	1.865	17	21	35	rVB2	32985	51275	3.69%	0.401%
2	2.225	75	80	82	rBV2	12220	18117	1.31%	0.142%
3	2.389	103	107	115	rVB	8181	14271	1.03%	0.112%
4	2.914	185	193	205	rBV	29536	65972	4.75%	0.516%
5	3.261	243	250	262	rVB4	6598	19846	1.43%	0.155%
6	3.956	350	364	378	rBV3	24789	83299	6.00%	0.652%
7	4.206	395	405	418	rBV	22025	61816	4.45%	0.484%
8	4.718	476	489	514	rBV4	55182	228046	16.43%	1.785%
9	5.242	559	575	593	rVB3	32859	127001	9.15%	0.994%
10	5.596	622	633	644	rVB4	5075	17876	1.29%	0.140%
11	5.755	649	659	668	rBV4	7184	21722	1.57%	0.170%
12	6.699	803	814	823	rBV2	24200	76296	5.50%	0.597%
13	6.797	824	830	839	rVB2	6684	17408	1.25%	0.136%
14	6.992	852	862	875	rVB2	14512	43186	3.11%	0.338%
15	7.529	943	950	961	rVB6	6050	18361	1.32%	0.144%
16	7.724	969	982	988	rBV	187751	458767	33.05%	3.592%
17	7.797	988	994	1002	rVV	323984	735625	53.00%	5.759%
18	7.876	1002	1007	1018	rVB	52338	125274	9.03%	0.981%
19	8.004	1018	1028	1035	rBV3	12003	30216	2.18%	0.237%
20	8.150	1042	1052	1063	rVB2	188050	454399	32.74%	3.557%
21	8.333	1063	1082	1092	rVV	338590	905221	65.22%	7.087%
22	8.419	1092	1096	1105	rVB5	8072	20120	1.45%	0.158%
23	8.693	1132	1141	1151	rBV	480846	961923	69.31%	7.531%
24	8.931	1172	1180	1186	rBV2	19680	43594	3.14%	0.341%
25	9.193	1208	1223	1227	rBV2	65207	160955	11.60%	1.260%
26	9.242	1227	1231	1238	rVV	52732	99972	7.20%	0.783%
27	9.327	1238	1245	1250	rVV	34116	69079	4.98%	0.541%
28	9.376	1250	1253	1260	rVV5	10078	21792	1.57%	0.171%
29	9.461	1260	1267	1274	rVB5	11079	28809	2.08%	0.226%
30	9.620	1285	1293	1302	rVB	312383	609037	43.88%	4.768%
31	9.748	1302	1314	1323	rBV	400857	844513	60.85%	6.611%
32	9.857	1327	1332	1339	rVB4	18083	37892	2.73%	0.297%
33	9.937	1339	1345	1351	rBV5	18697	48568	3.50%	0.380%
34	9.992	1351	1354	1362	rVB4	12135	25303	1.82%	0.198%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	10.113	1366	1374	1379	rBV	148811	262567	18.92%	2.056%
36	10.181	1379	1385	1397	rVB	793028	1387952	100.00%	10.866%
37	10.467	1428	1432	1437	rBV3	10713	18709	1.35%	0.146%
38	10.522	1437	1441	1445	rBV2	9261	14919	1.07%	0.117%
39	10.607	1445	1455	1457	rBV4	25585	67458	4.86%	0.528%
40	10.711	1466	1472	1477	rVB4	11150	22413	1.61%	0.175%
41	10.802	1477	1487	1492	rBV3	9823	21655	1.56%	0.170%
42	10.900	1492	1503	1514	rVV4	15102	57037	4.11%	0.447%
43	11.004	1516	1520	1524	rVV3	11450	19002	1.37%	0.149%
44	11.089	1528	1534	1539	rVB4	12093	22650	1.63%	0.177%
45	11.254	1556	1561	1566	rBV6	10565	24296	1.75%	0.190%
46	11.314	1567	1571	1577	rVB5	17474	34903	2.51%	0.273%
47	11.491	1592	1600	1606	rBV	735179	1207577	87.00%	9.454%
48	11.741	1635	1641	1645	rBV5	11480	22320	1.61%	0.175%
49	11.997	1678	1683	1691	rBV7	7600	19867	1.43%	0.156%
50	12.326	1731	1737	1743	rBV	107975	161589	11.64%	1.265%
51	12.479	1752	1762	1770	rBV	624985	1008761	72.68%	7.897%
52	13.247	1882	1888	1895	rVV3	16100	40294	2.90%	0.315%
53	13.424	1910	1917	1925	rBV	788776	1199256	86.40%	9.389%
54	13.521	1929	1933	1939	rBV2	10820	17294	1.25%	0.135%
55	13.588	1939	1944	1947	rBV	47020	72694	5.24%	0.569%
56	13.625	1947	1950	1954	rVV	51087	74227	5.35%	0.581%
57	13.674	1954	1958	1965	rVV2	23815	41787	3.01%	0.327%
58	13.753	1966	1971	1976	rVB4	14061	23151	1.67%	0.181%
59	13.966	2000	2006	2011	rBV3	8230	15218	1.10%	0.119%
60	14.027	2011	2016	2020	rVV	9442	16470	1.19%	0.129%
61	14.082	2020	2025	2030	rVB	28192	46539	3.35%	0.364%
62	14.296	2056	2060	2064	rVB2	9449	15528	1.12%	0.122%
63	14.375	2068	2073	2077	rBV3	10701	17746	1.28%	0.139%
64	14.454	2082	2086	2093	rBV2	27861	41949	3.02%	0.328%
65	14.564	2100	2104	2109	rBV3	25236	41131	2.96%	0.322%
66	14.661	2115	2120	2121	rVV	39357	49523	3.57%	0.388%
67	14.686	2121	2124	2131	rVB	49063	79384	5.72%	0.621%
68	15.058	2178	2185	2191	rVB3	14699	31311	2.26%	0.245%
69	15.228	2203	2213	2217	rBV3	5768	15620	1.13%	0.122%
70	15.692	2282	2289	2297	rVB5	5415	15115	1.09%	0.118%

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Instrument :
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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

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Title : SW846 8260

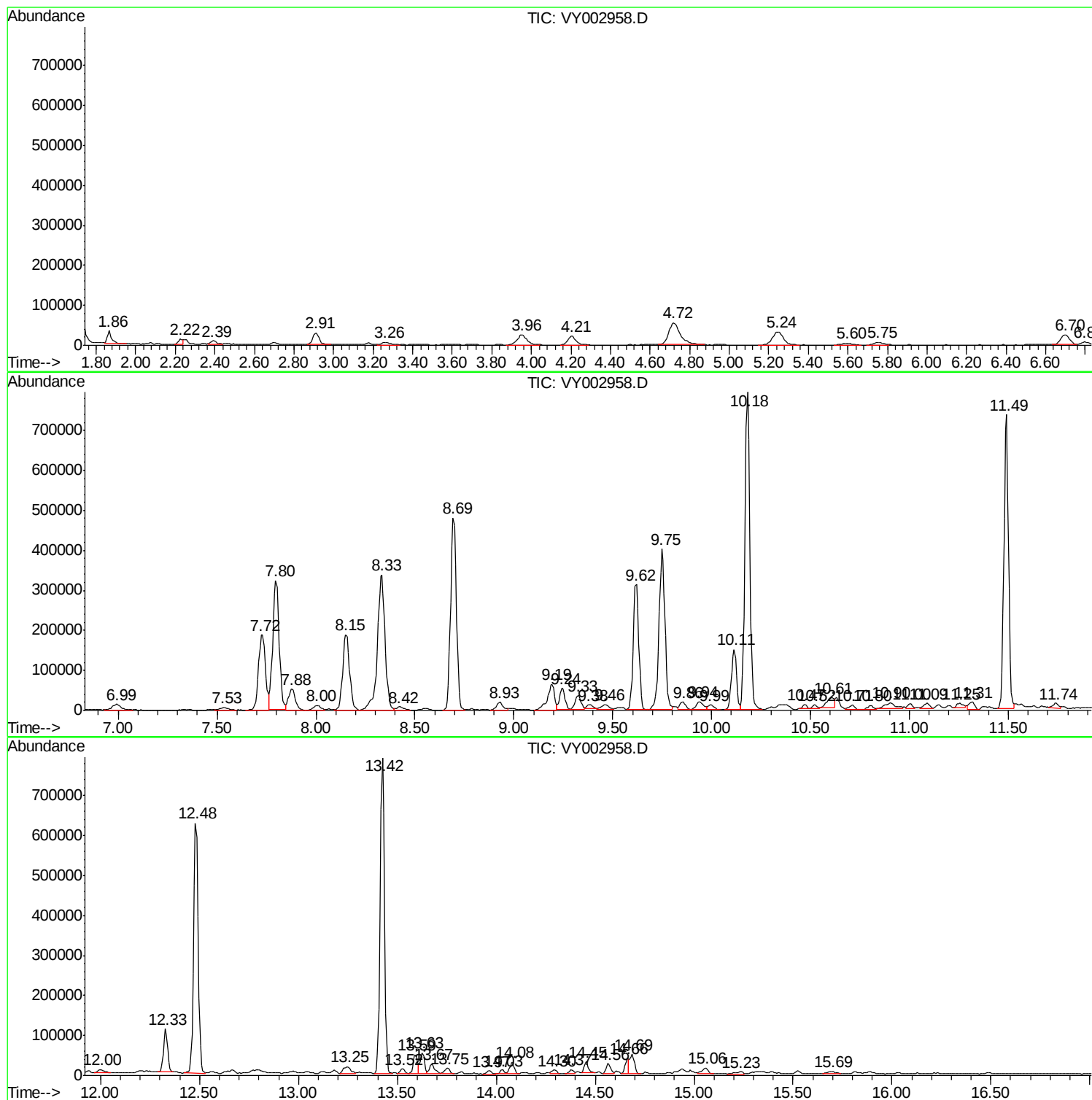
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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



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ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TW-11

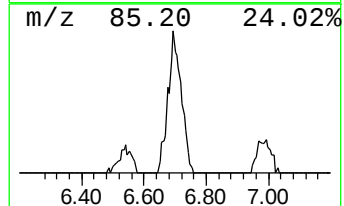
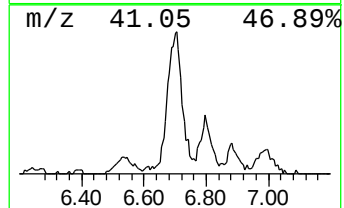
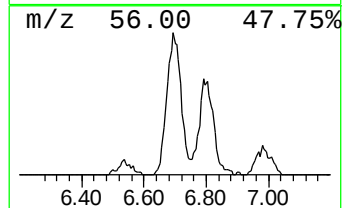
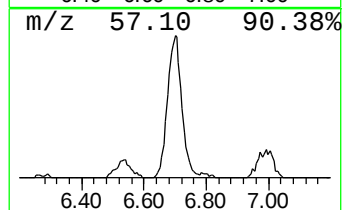
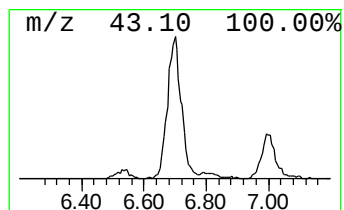
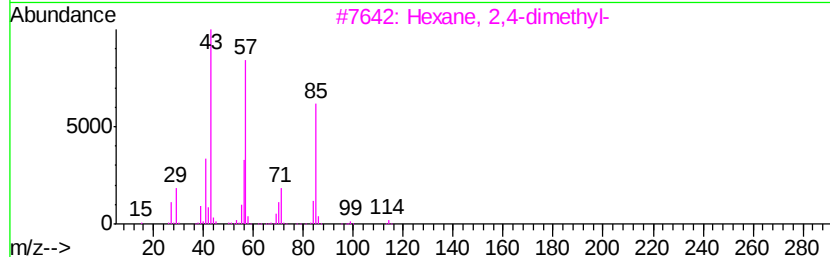
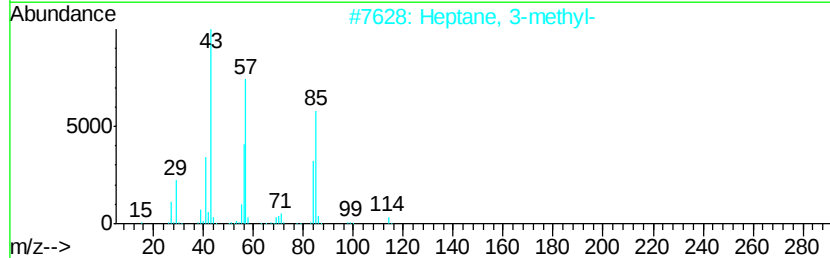
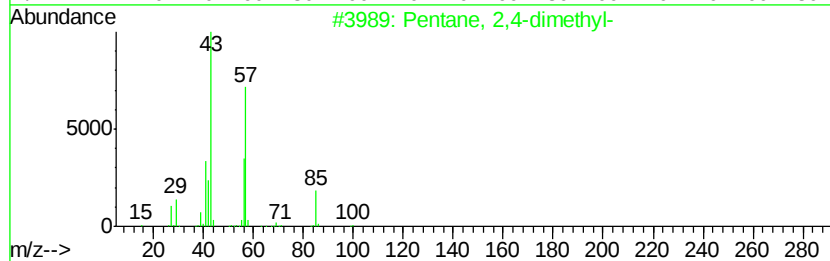
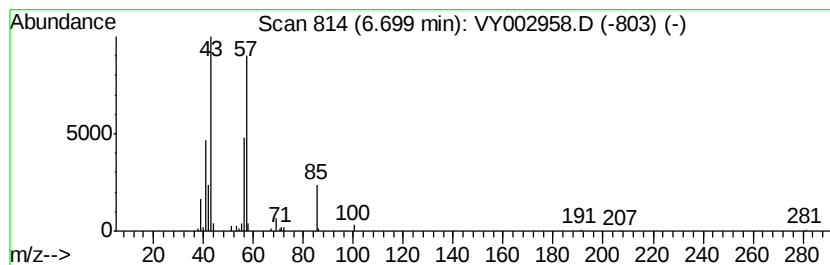
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 2 Pentane, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	5.19 ug/l	76296	Pentafluorobenzene	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	78
2		Heptane, 3-methyl-	114	C8H18	000589-81-1	64
3		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4		.alpha.-Amino-.gamma.-butyrolactone	101	C4H7NO2	001192-20-7	59
5		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	53



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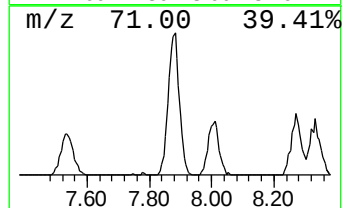
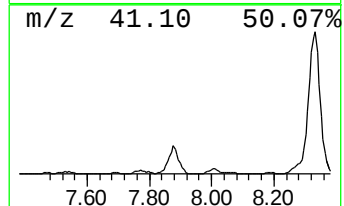
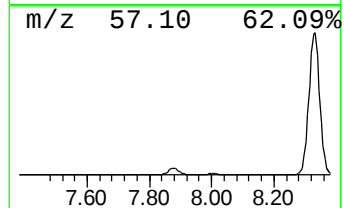
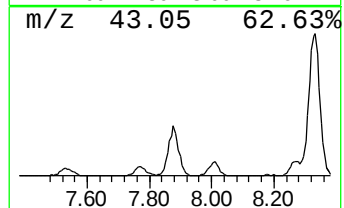
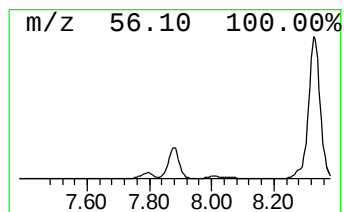
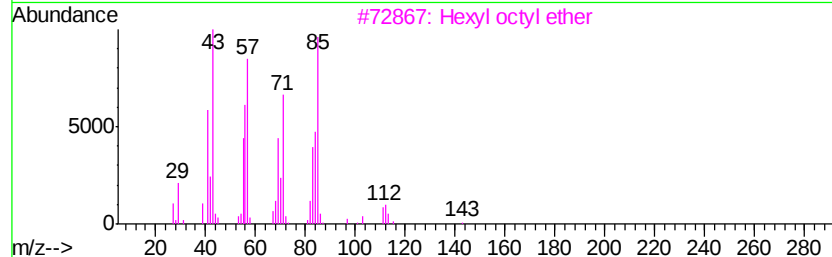
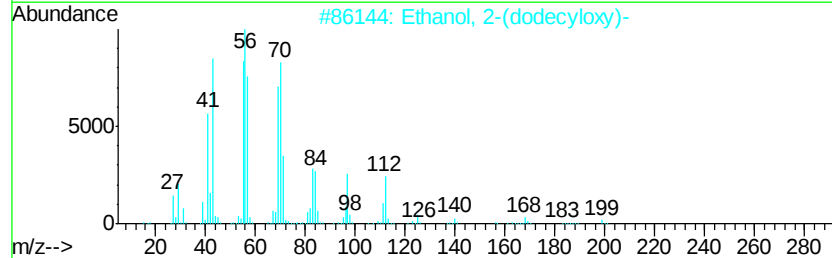
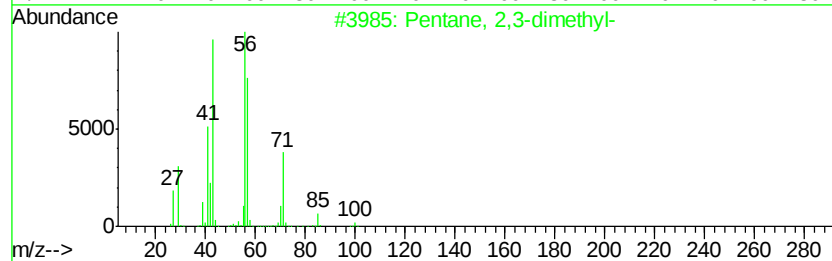
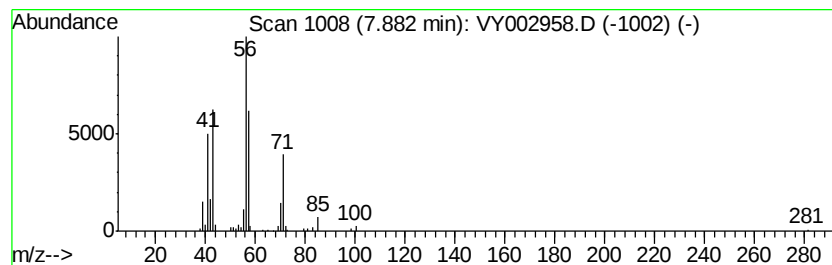
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	8.51 ug/l	125274	Pentafluorobenzene	7.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	64
3			Hexyl octyl ether	214	C14H30O	017071-54-4	47
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45
5			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	38



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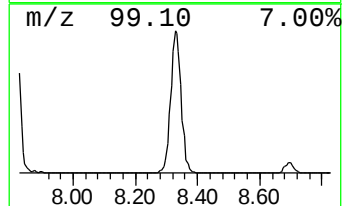
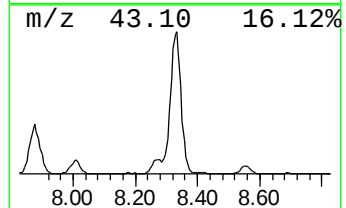
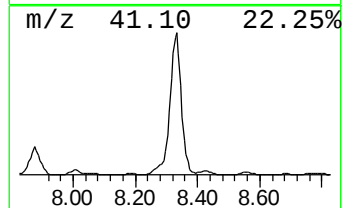
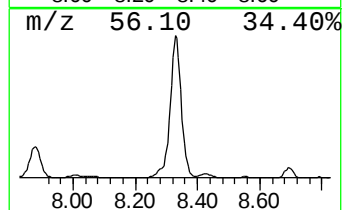
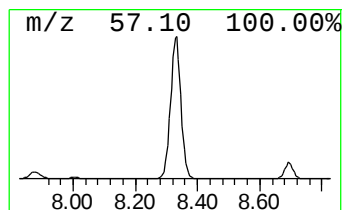
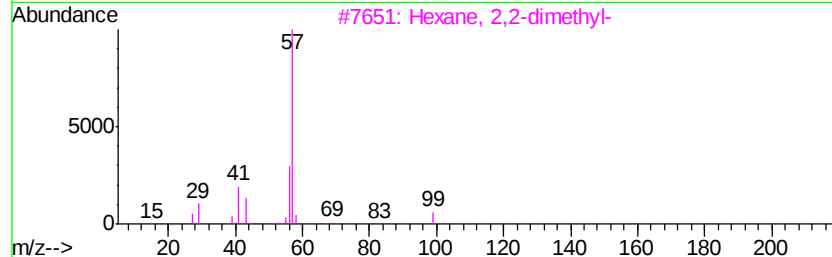
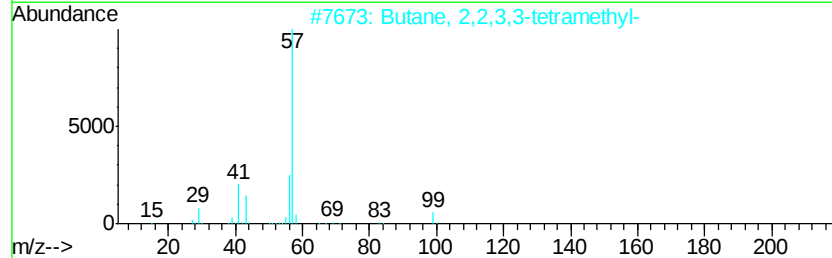
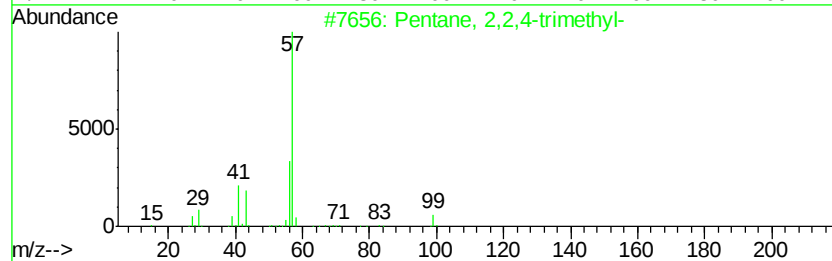
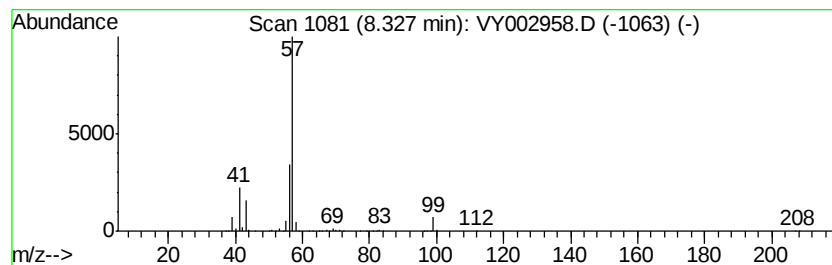
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentane, 2,2,4-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.33	47.05 ug/l	905221	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45



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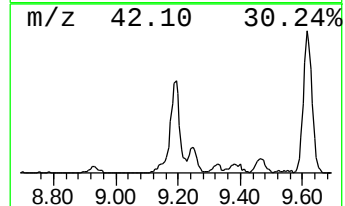
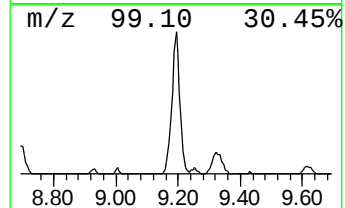
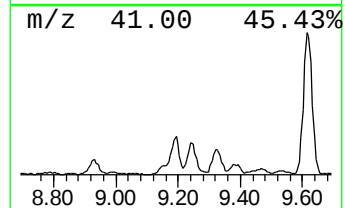
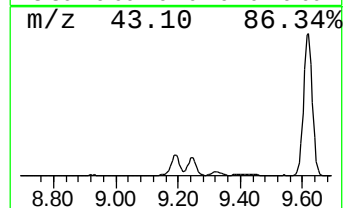
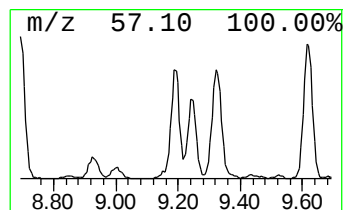
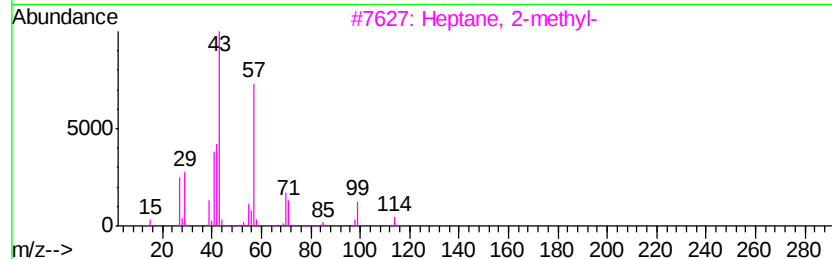
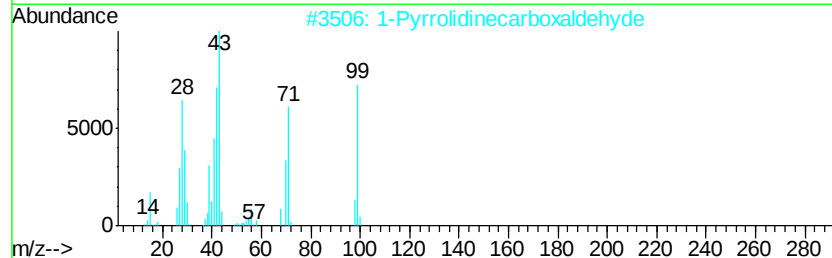
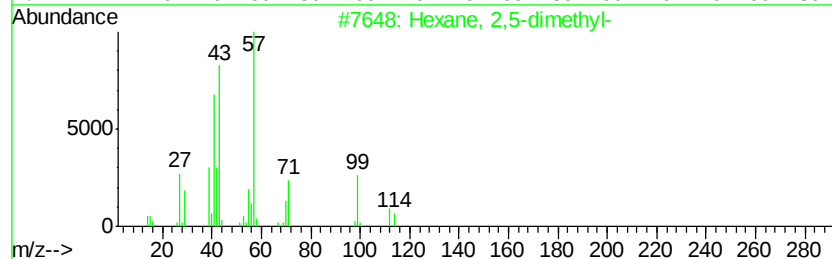
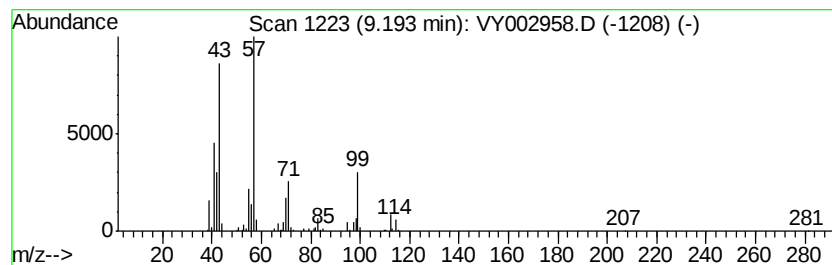
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexane, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.19	8.37 ug/l	160955	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	87
2		1-Pyrrolidinecarboxaldehyde	99	C5H9NO	003760-54-1	52
3		Heptane, 2-methyl-	114	C8H18	000592-27-8	47
4		Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	38
5		2-Piperidinone	99	C5H9NO	000675-20-7	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY061920\
Data File : VY002958.D
Acq On : 19 Jun 2020 23:22
Operator : SY/MD
Sample : L3063-15
Misc : 5.06G/5ML/MSVOA Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TW-11

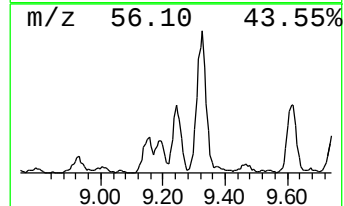
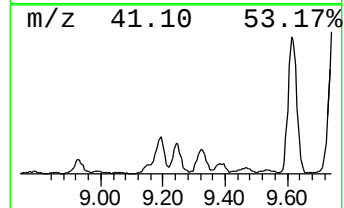
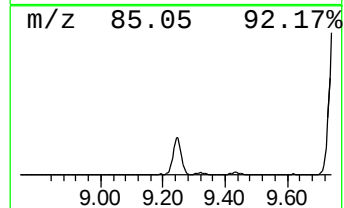
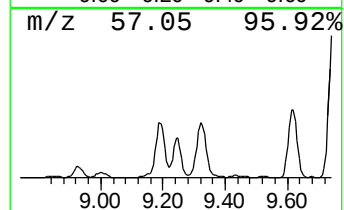
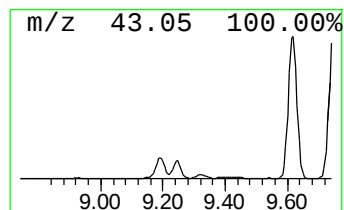
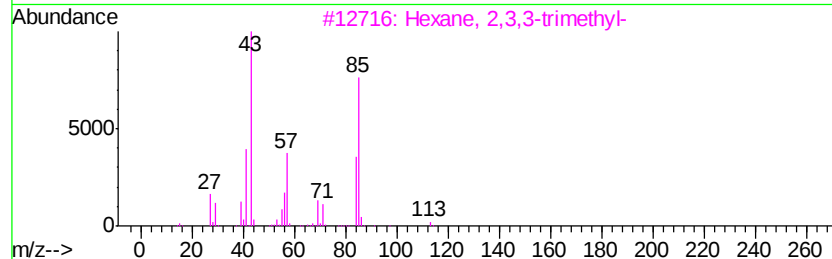
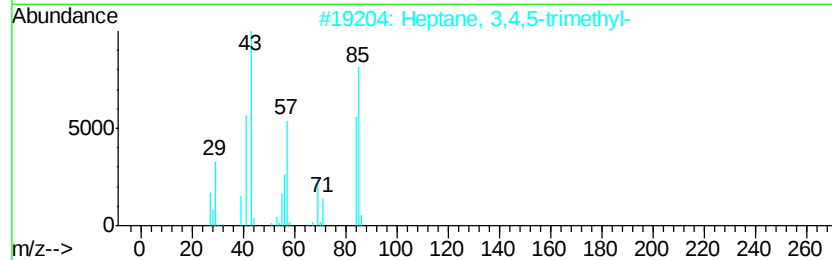
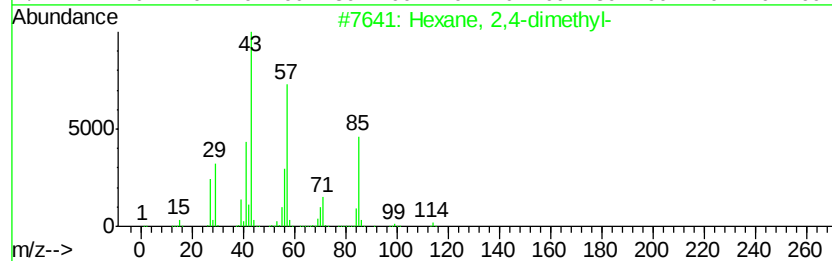
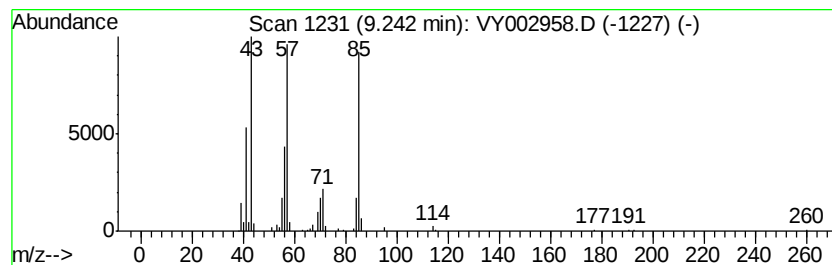
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Quant Title : SW846 8260

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

Peak Number 6 Hexane, 2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	5.20 ug/l	99972	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
2		Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	72
3		Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	59
4		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59
5		4,4-Dimethyl octane	142	C10H22	015869-95-1	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY061920\
Data File : VY002958.D
Acq On : 19 Jun 2020 23:22
Operator : SY/MD
Sample : L3063-15
Misc : 5.06G/5ML/MSVOA Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TW-11

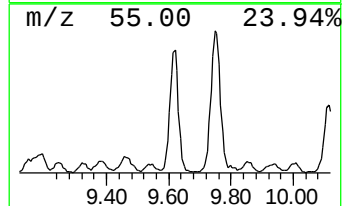
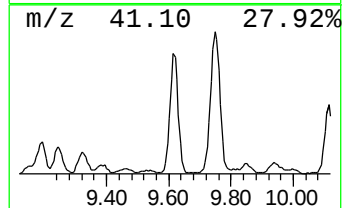
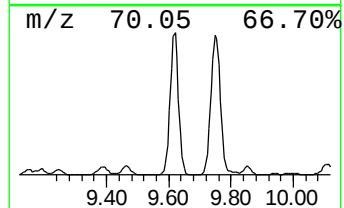
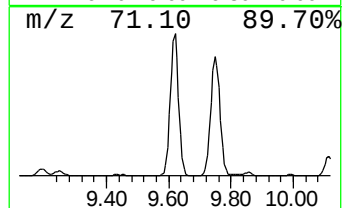
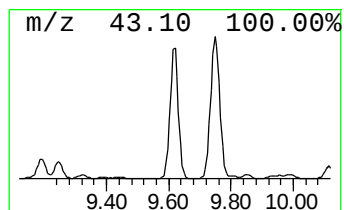
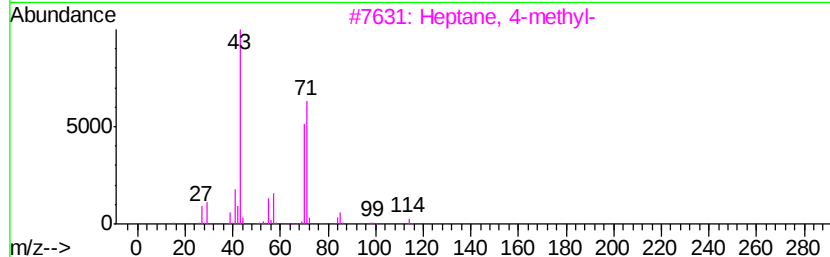
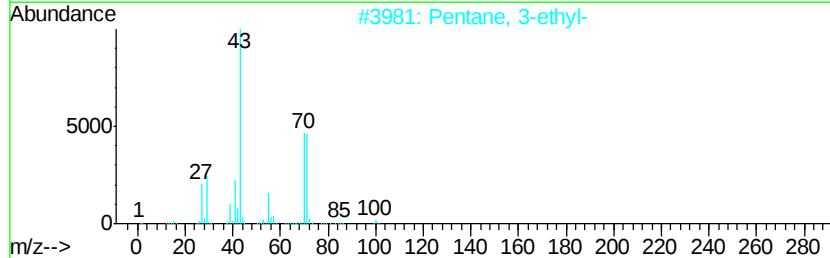
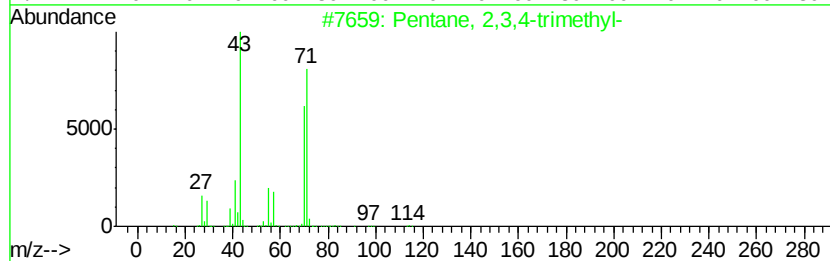
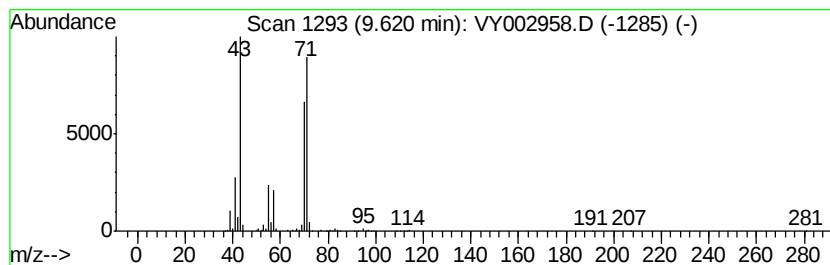
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Quant Title : SW846 8260

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

Peak Number 7 Pentane, 2,3,4-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	31.66 ug/l	609037	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4		Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	78
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY061920\
Data File : VY002958.D
Acq On : 19 Jun 2020 23:22
Operator : SY/MD
Sample : L3063-15
Misc : 5.06G/5ML/MSVOA Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TW-11

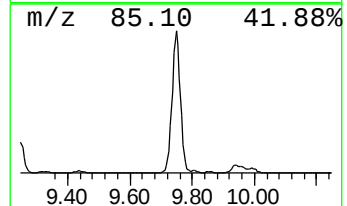
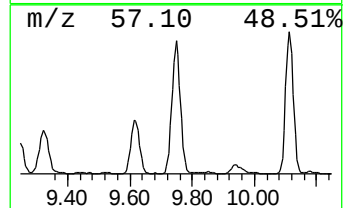
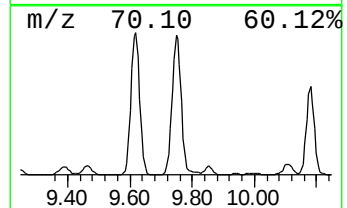
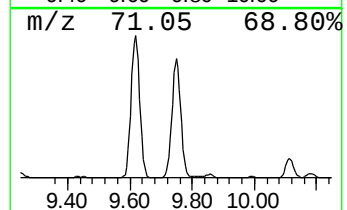
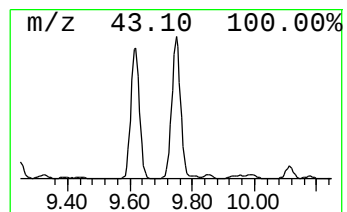
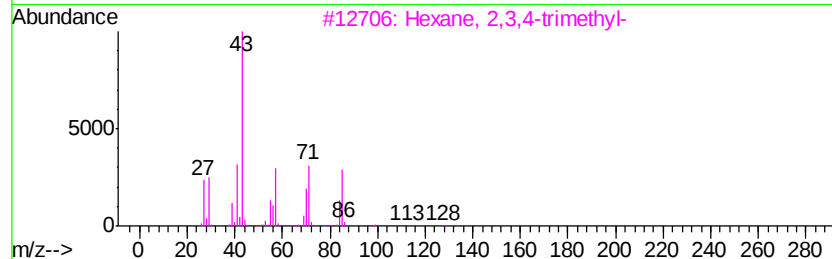
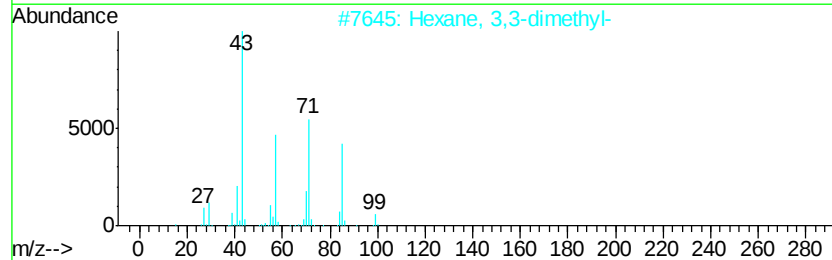
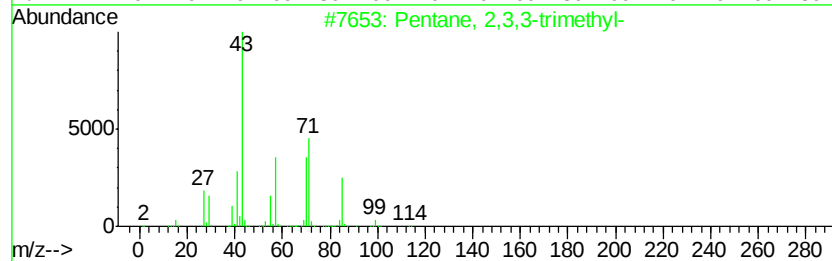
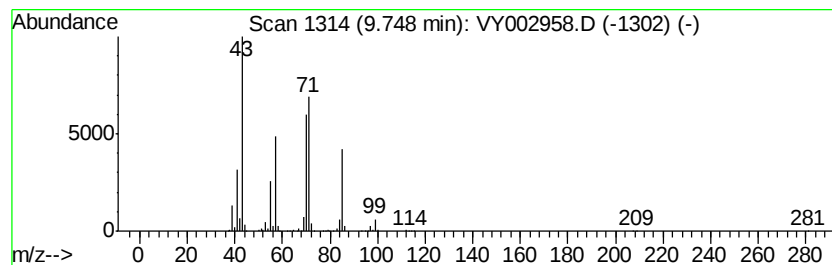
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 8 Pentane, 2,3,3-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.75	43.90 ug/l	844513	1,4-Difluorobenzene	8.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
3		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	53
5		Heptane, 4-methyl-	114	C8H18	000589-53-7	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY061920\
Data File : VY002958.D
Acq On : 19 Jun 2020 23:22
Operator : SY/MD
Sample : L3063-15
Misc : 5.06G/5ML/MSVOA Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TW-11

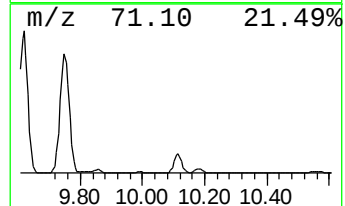
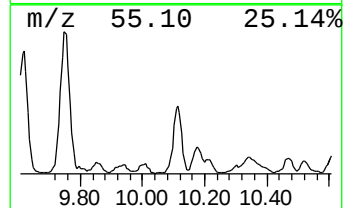
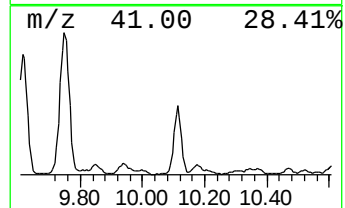
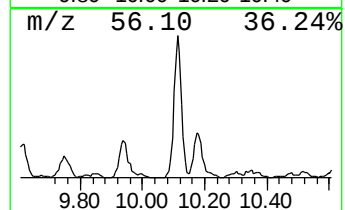
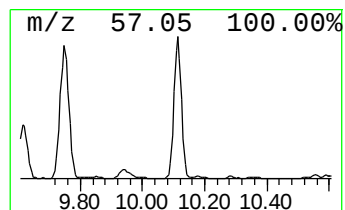
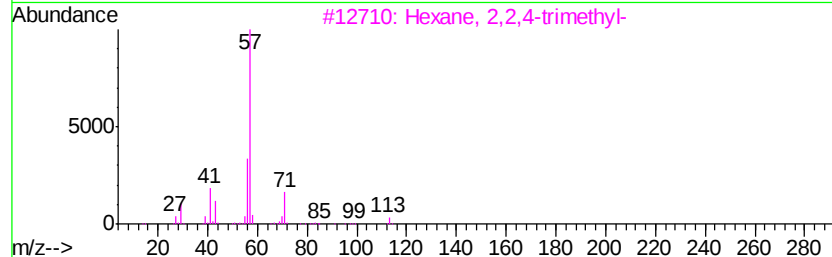
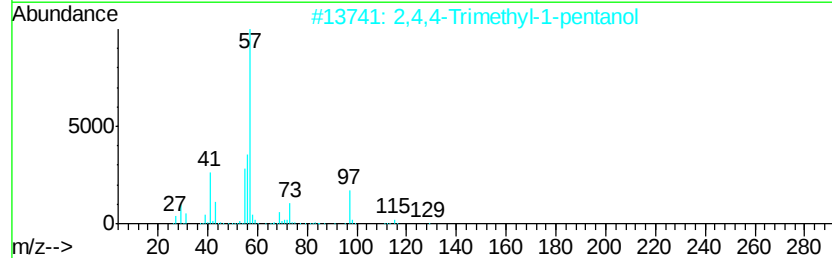
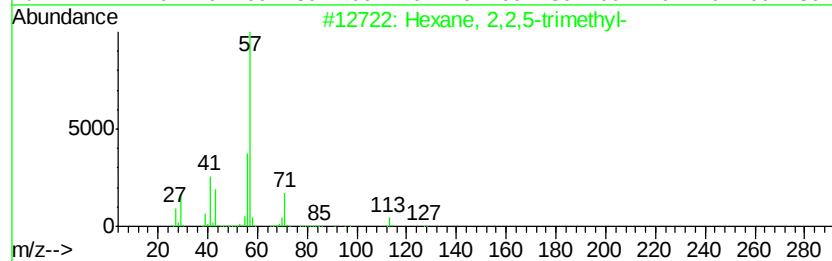
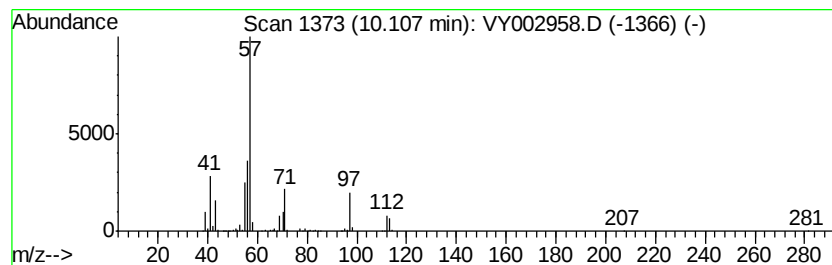
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y061920S.M
Quant Title : SW846 8260

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

Peak Number 9 Hexane, 2,2,5-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.11	10.87 ug/l	262567	Chlorobenzene-d5	11.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	50	
2	2,4,4-Trimethyl-1-pentanol	130	C8H18O	016325-63-6	50	
3	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	50	
4	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	47	
5	1-Hexene, 5,5-dimethyl-	112	C8H16	007116-86-1	47	



Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_Y\DATA\VY061920\
Data File : VY002958.D
Acq On : 19 Jun 2020 23:22
Operator : SY/MD
Sample : L3063-15
Misc : 5.06G/5ML/MSV0A_Y/SOIL
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSV0A_Y
ClientSampleId :
TW-11

Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_Y\METHODS\82Y061920S.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
Pentane, 2,4-dime...	6.70	5.2	ug/l	76296	1	7.80	735625	50.0
Pentane, 2,3-dime...	7.88	8.5	ug/l	125274	1	7.80	735625	50.0
Pentane, 2,2,4-tr...	8.33	47.0	ug/l	905221	2	8.69	961923	50.0
Hexane, 2,5-dimet...	9.19	8.4	ug/l	160955	2	8.69	961923	50.0
Hexane, 2,4-dimet...	9.24	5.2	ug/l	99972	2	8.69	961923	50.0
Pentane, 2,3,4-tr...	9.62	31.7	ug/l	609037	2	8.69	961923	50.0
Pentane, 2,3,3-tr...	9.75	43.9	ug/l	844513	2	8.69	961923	50.0
Hexane, 2,2,5-tri...	10.11	10.9	ug/l	262567	3	11.49	1207580	50.0