

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
 Data File : VY009602.D
 Acq On : 19 Jul 2022 20:10
 Operator : KP/MD
 Sample : N3652-05
 Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SED-3-(6-12)

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

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

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
 Title : SW846 8260

Signal : TIC: VY009602.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.521	914	935	951	rBV	258270	785538	1.28%	0.215%
2	7.783	972	978	985	rVB	315429	691750	1.13%	0.190%
3	8.319	1048	1066	1090	rBV	22048090	60093570	98.19%	16.465%
4	8.685	1116	1126	1133	rBV	526437	1075876	1.76%	0.295%
5	9.319	1222	1230	1237	rBV	3048910	6162009	10.07%	1.688%
6	9.459	1244	1253	1261	rBV2	561781	1350357	2.21%	0.370%
7	9.618	1270	1279	1290	rVB	24218245	51481925	84.12%	14.106%
8	9.746	1290	1300	1312	rBV	27723365	61198483	100.00%	16.768%
9	9.856	1312	1318	1324	rVB3	341941	813321	1.33%	0.223%
10	9.929	1324	1330	1336	rBV4	574308	1147839	1.88%	0.314%
11	9.996	1336	1341	1347	rVB3	517172	1081653	1.77%	0.296%
12	10.112	1347	1360	1366	rBV	11438812	20694537	33.82%	5.670%
13	10.179	1366	1371	1383	rVB2	818291	1863538	3.05%	0.511%
14	10.350	1391	1399	1409	rVB3	530224	1857136	3.03%	0.509%
15	10.478	1409	1420	1425	rBV	1516174	2848531	4.65%	0.780%
16	10.636	1437	1446	1455	rBV2	2856520	6410688	10.48%	1.756%
17	10.764	1462	1467	1470	rBV3	362087	783341	1.28%	0.215%
18	10.807	1470	1474	1479	rVV	984190	1693997	2.77%	0.464%
19	10.892	1479	1488	1496	rVV5	864502	3171886	5.18%	0.869%
20	11.002	1502	1506	1511	rVB2	1241047	1970416	3.22%	0.540%
21	11.087	1511	1520	1525	rBV4	2591928	7322839	11.97%	2.006%
22	11.142	1525	1529	1541	rVB4	2752962	6065838	9.91%	1.662%
23	11.246	1541	1546	1552	rBV3	1187515	3020822	4.94%	0.828%
24	11.313	1552	1557	1563	rVB3	1634642	3930083	6.42%	1.077%
25	11.410	1563	1573	1580	rBV6	1578726	5585298	9.13%	1.530%
26	11.483	1580	1585	1590	rVV2	1904710	4566231	7.46%	1.251%
27	11.532	1590	1593	1594	rVV2	836635	1099061	1.80%	0.301%
28	11.563	1594	1598	1602	rVB	2465000	3758412	6.14%	1.030%
29	11.624	1602	1608	1613	rBV3	1293623	2199854	3.59%	0.603%
30	11.672	1613	1616	1620	rVB4	638716	891855	1.46%	0.244%
31	11.740	1620	1627	1637	rVB4	3187352	7569301	12.37%	2.074%
32	11.837	1637	1643	1646	rBV4	704069	1490088	2.43%	0.408%
33	11.874	1646	1649	1651	rBV	428848	614519	1.00%	0.168%
34	11.935	1656	1659	1663	rBV	1440843	2318396	3.79%	0.635%
35	11.989	1663	1668	1675	rVB5	1177697	2992674	4.89%	0.820%
36	12.057	1675	1679	1681	rBV3	1070874	1479539	2.42%	0.405%

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Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
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37	12.087	1681	1684	1691	rVB2	1494659	3313821	5.41%	0.908%
38	12.160	1691	1696	1697	rBV2	484393	702899	1.15%	0.193%
39	12.203	1698	1703	1705	rBV4	1184979	2218512	3.63%	0.608%
40	12.459	1740	1745	1754	rVB	6536730	11409559	18.64%	3.126%
41	12.569	1754	1763	1768	rBV2	2156041	5167986	8.44%	1.416%
42	12.624	1768	1772	1782	rVB	2000417	5407712	8.84%	1.482%
43	12.715	1782	1787	1793	rBV8	943261	2605882	4.26%	0.714%
44	12.947	1821	1825	1829	rVB4	927459	1216799	1.99%	0.333%
45	13.008	1829	1835	1838	rBV8	463103	904570	1.48%	0.248%
46	13.069	1840	1845	1853	rVB	1670144	3331822	5.44%	0.913%
47	13.166	1857	1861	1867	rVB	2280852	3830502	6.26%	1.050%
48	13.245	1867	1874	1876	rBV4	719650	1735451	2.84%	0.475%
49	13.282	1876	1880	1888	rVB3	2830438	4420058	7.22%	1.211%
50	13.422	1895	1903	1907	rBV4	712914	2008994	3.28%	0.550%
51	13.483	1908	1913	1922	rVB	1912538	4139499	6.76%	1.134%
52	13.562	1922	1926	1929	rBV4	455689	646861	1.06%	0.177%
53	13.642	1936	1939	1943	rVB	730338	880626	1.44%	0.241%
54	13.812	1963	1967	1971	rBV3	606688	1170193	1.91%	0.321%
55	13.995	1995	1997	2002	rVB4	791535	1120646	1.83%	0.307%
56	14.074	2006	2010	2014	rBV	1252136	1611403	2.63%	0.442%
57	14.148	2015	2022	2028	rBV4	2157301	4875374	7.97%	1.336%
58	14.257	2034	2040	2046	rBV5	1508883	3424690	5.60%	0.938%
59	14.385	2055	2061	2063	rBV3	934278	1811295	2.96%	0.496%
60	14.471	2071	2075	2082	rVB5	750294	1425827	2.33%	0.391%
61	14.544	2084	2087	2090	rBV3	434292	744667	1.22%	0.204%
62	14.623	2096	2100	2104	rBV3	621613	1013074	1.66%	0.278%
63	14.727	2114	2117	2124	rVB6	796861	1547223	2.53%	0.424%
64	14.806	2126	2130	2133	rBV4	573970	938768	1.53%	0.257%
65	14.849	2134	2137	2139	rBV2	479967	680389	1.11%	0.186%
66	14.940	2149	2152	2155	rVB3	725563	968720	1.58%	0.265%
67	14.983	2156	2159	2165	rBV6	357964	897192	1.47%	0.246%
68	15.282	2203	2208	2213	rBV5	772737	1519773	2.48%	0.416%
69	15.446	2231	2235	2241	rBV5	619348	1355491	2.21%	0.371%
70	16.476	2400	2404	2419	rVB8	775343	2467054	4.03%	0.676%
71	16.763	2446	2451	2455	rBV3	707731	1381726	2.26%	0.379%

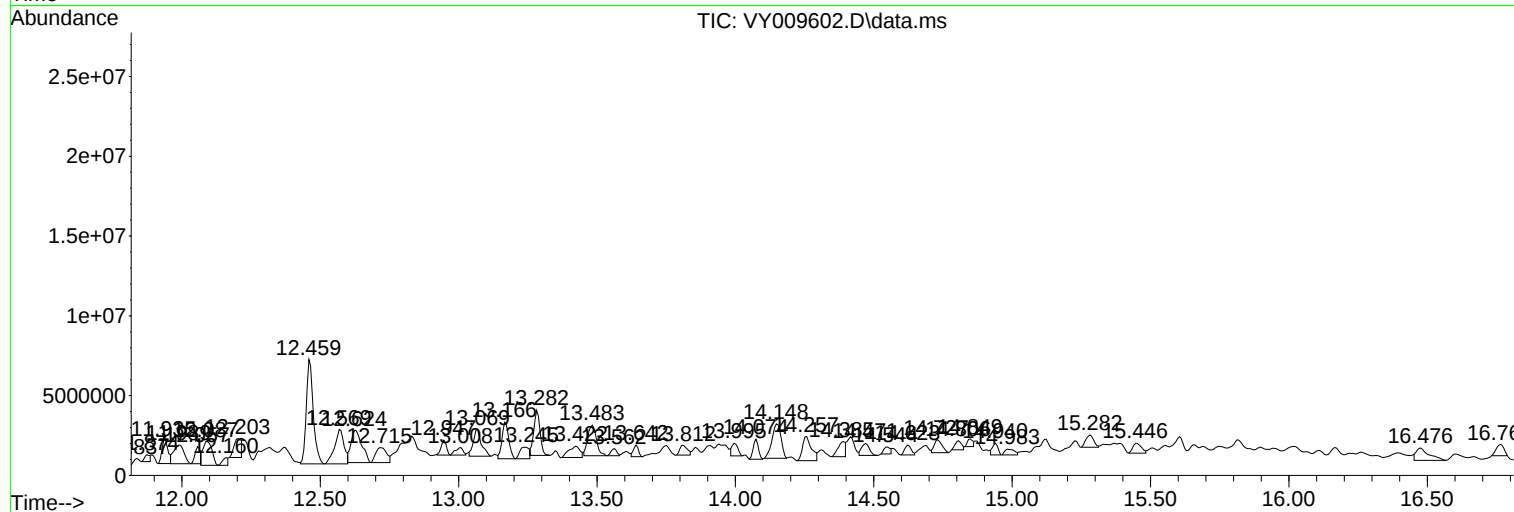
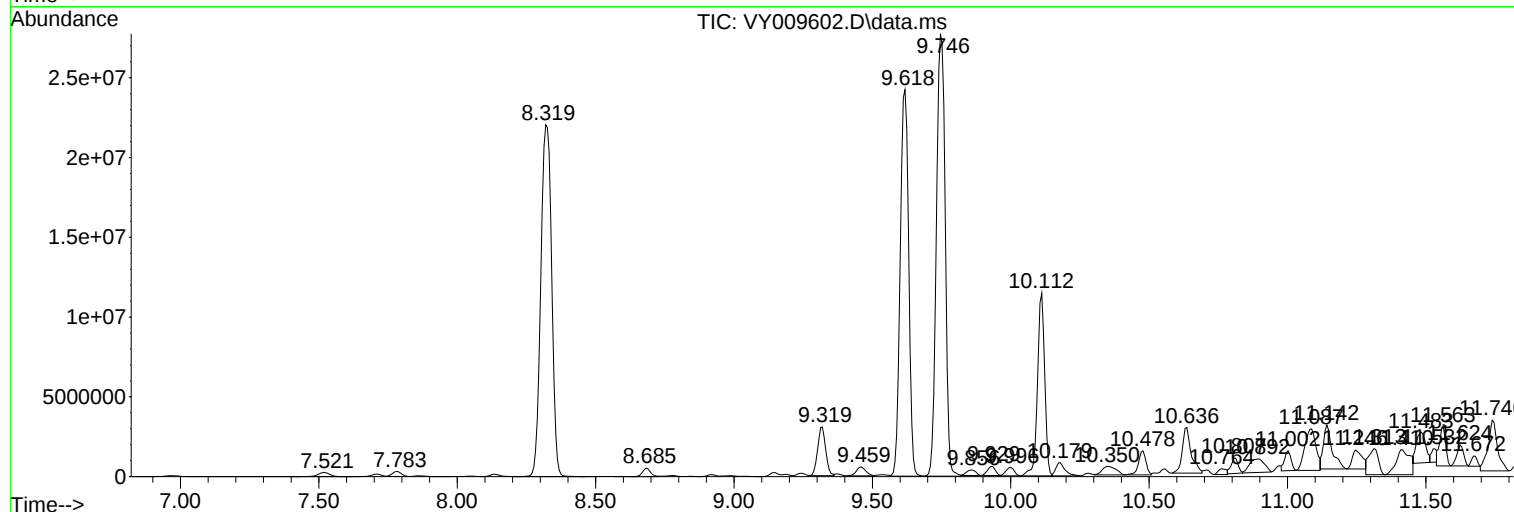
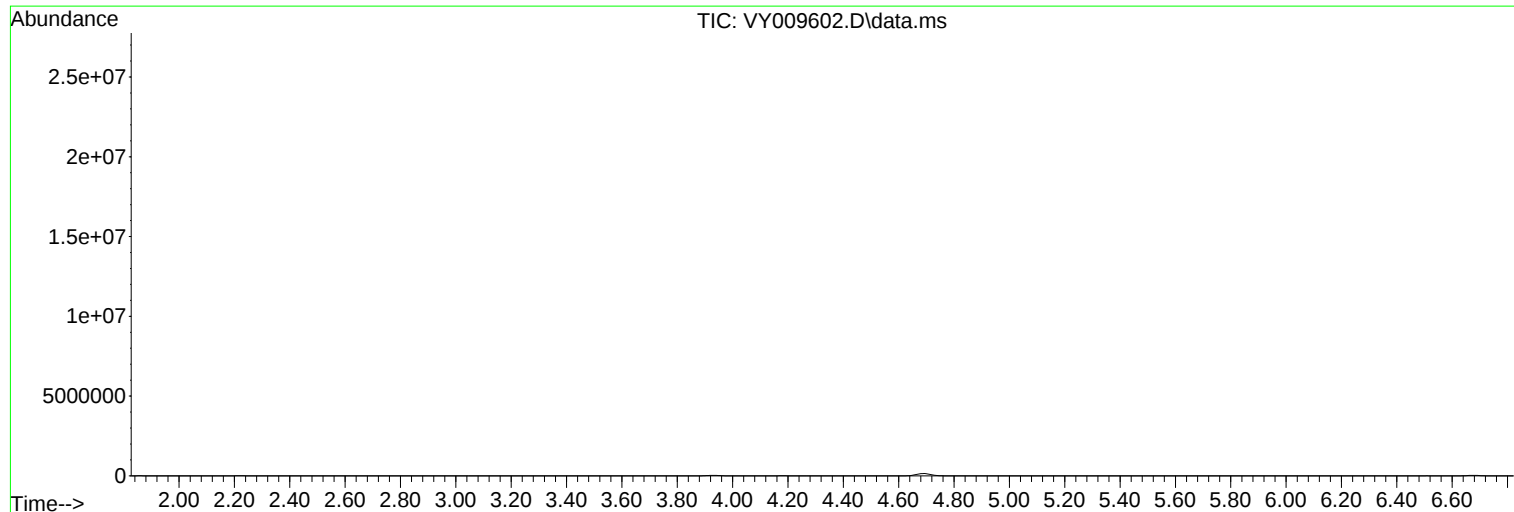
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Instrument :
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ClientSampleId :
SED-3-(6-12)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

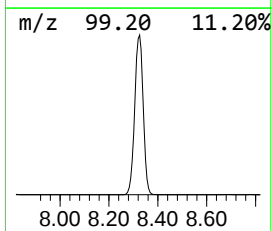
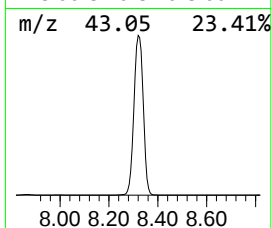
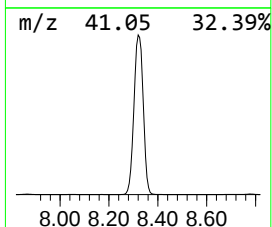
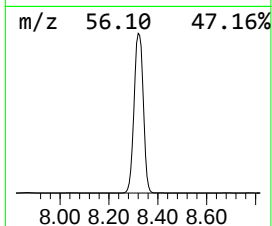
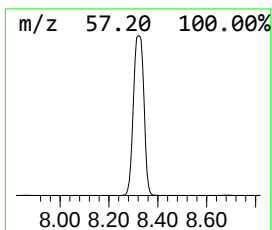
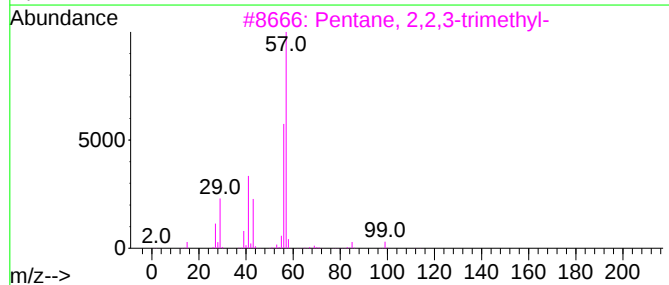
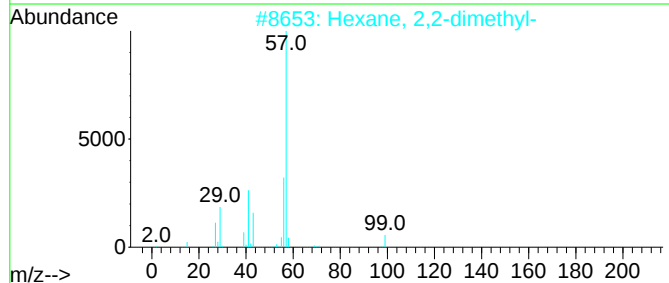
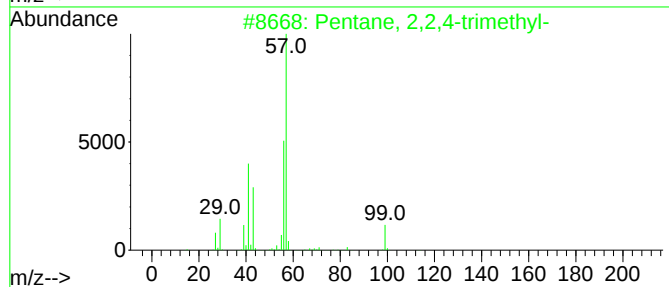
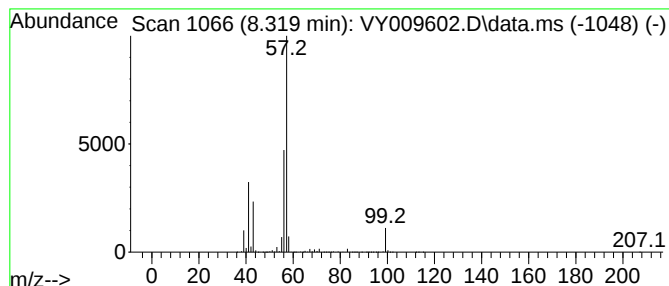
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.319	2792.77 ug/l	60093600	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
3			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	64
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59



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Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

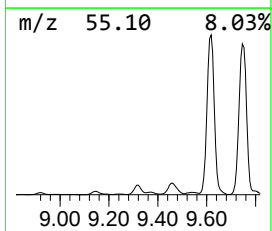
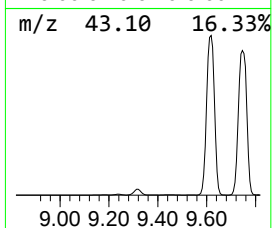
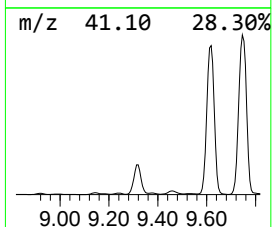
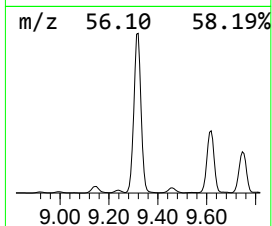
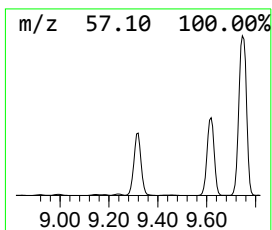
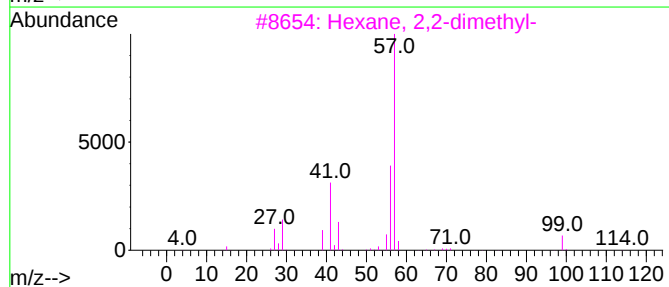
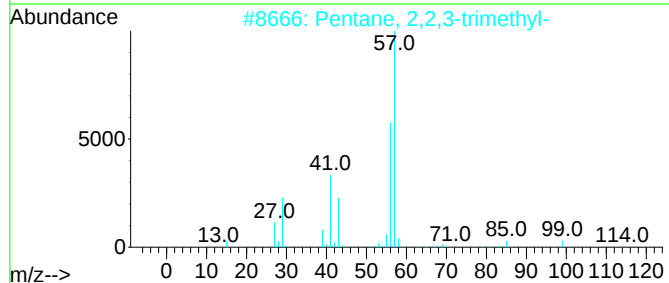
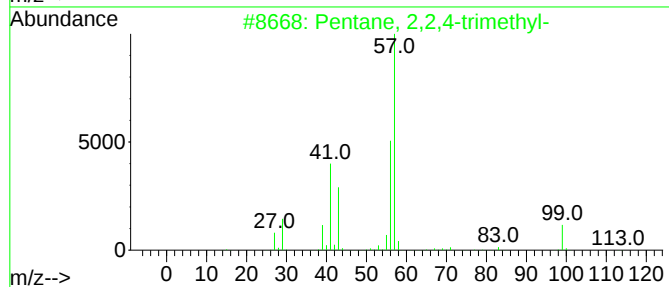
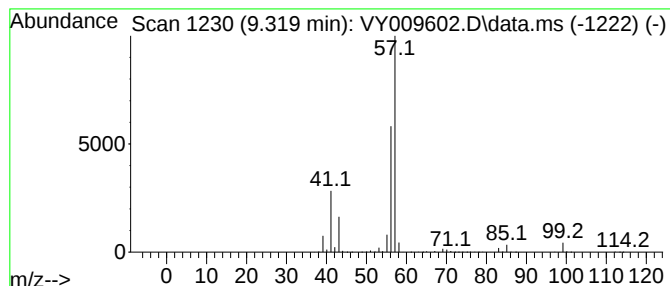
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.319	286.37 ug/l	6162010	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	72
3			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
5			Acetamide, N-2-propenyl-	99	C5H9NO	000692-33-1	50



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Instrument :
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ClientSampleId :
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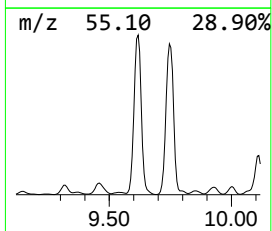
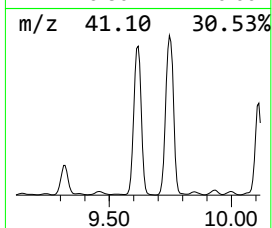
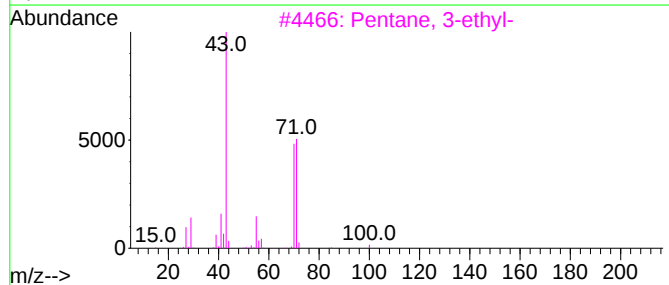
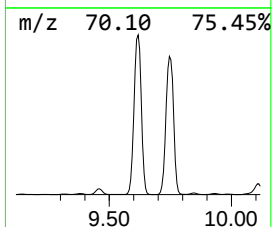
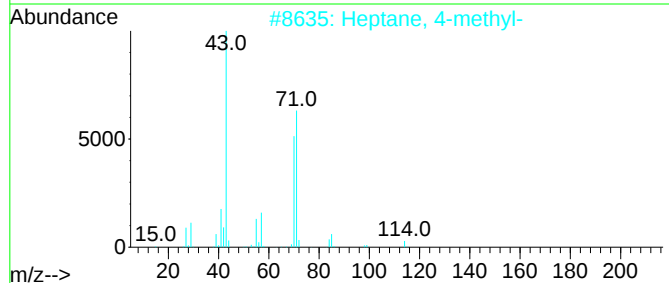
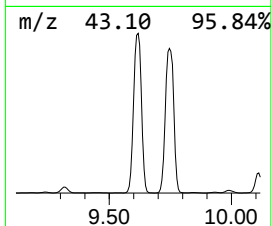
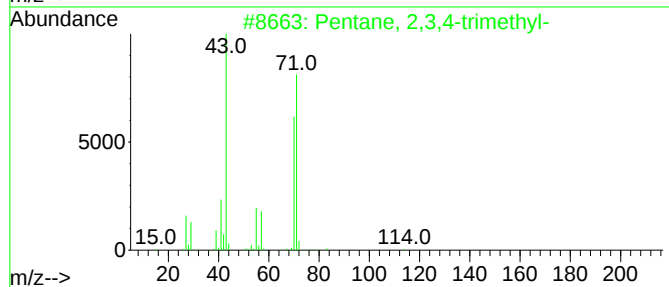
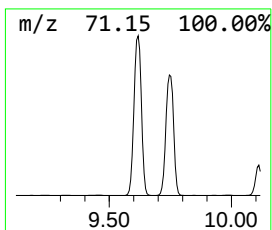
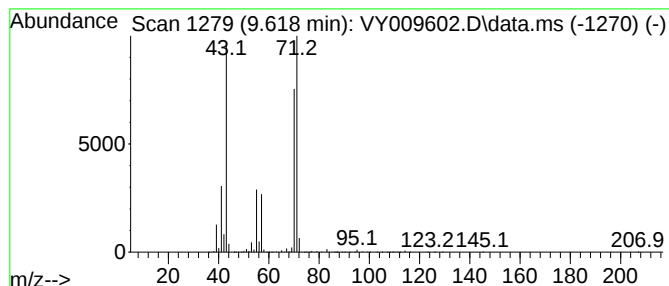
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.618	2392.56 ug/l	51481900	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	90
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	74
3			Pentane, 3-ethyl-	100	C7H16	000617-78-7	74
4			Diisooamyl ether	158	C10H22O	000544-01-4	74
5			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64



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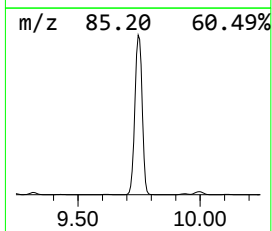
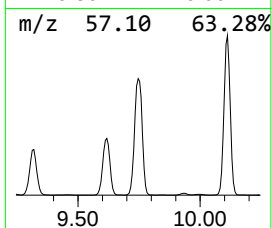
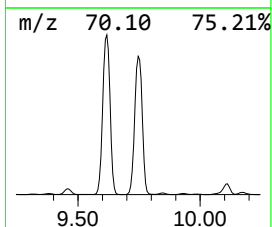
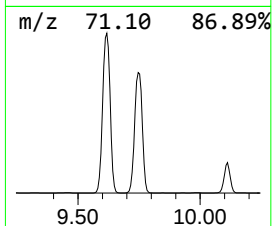
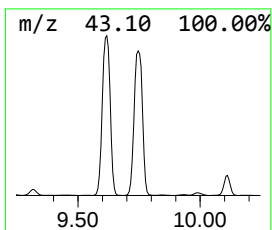
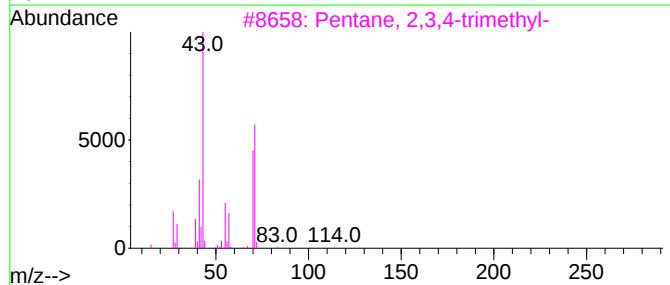
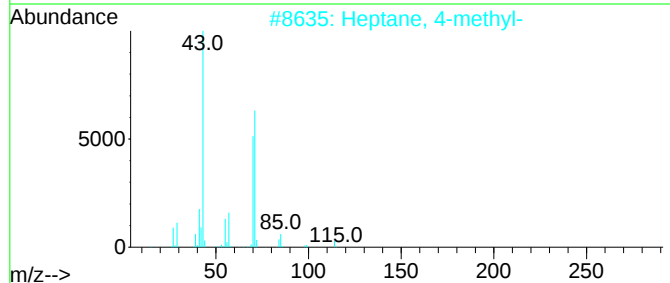
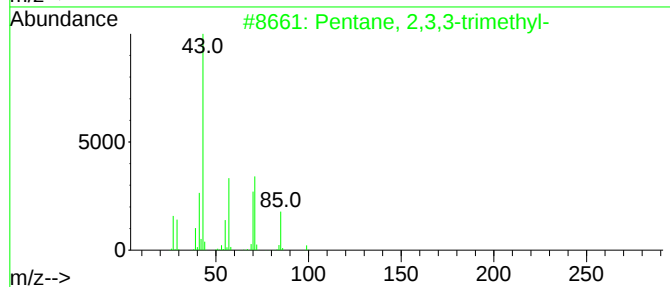
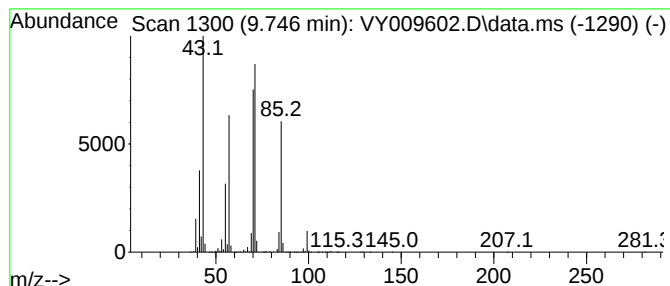
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.746	2844.12 ug/l	61198500	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
4			Pyrrolidine	71	C4H9N	000123-75-1	53
5			Diisooamyl ether	158	C10H22O	000544-01-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

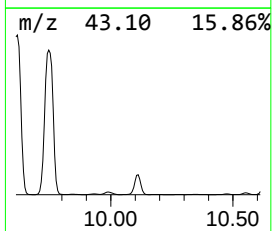
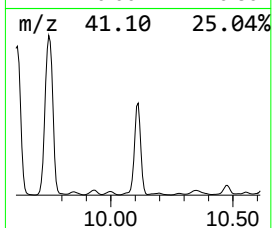
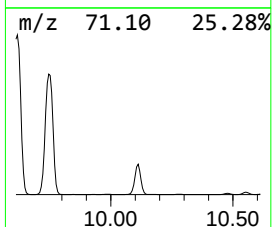
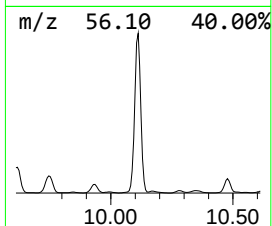
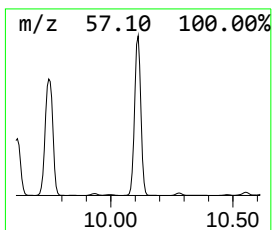
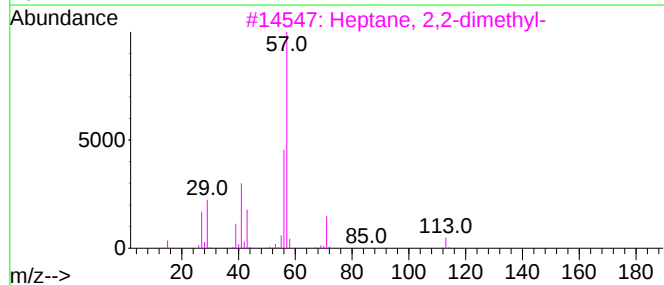
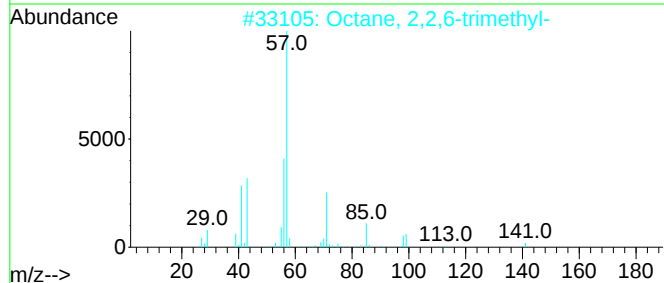
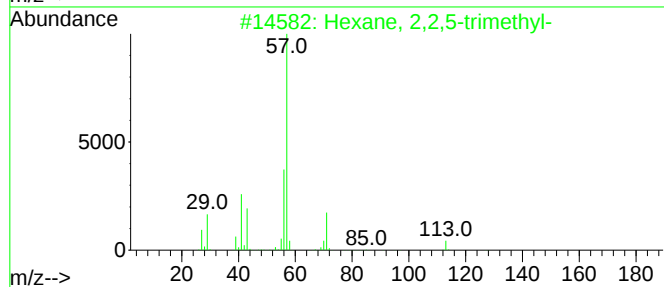
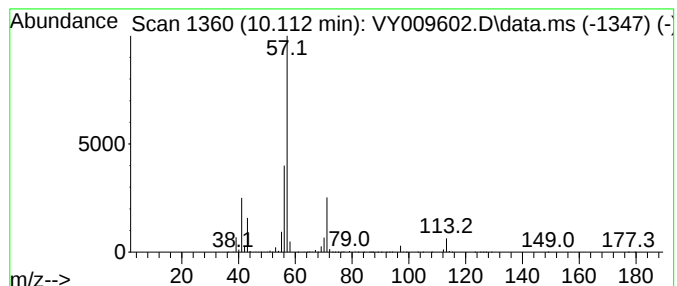
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.112	226.60 ug/l	20694500	Chlorobenzene-d5	11.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	83
2			Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	72
3			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

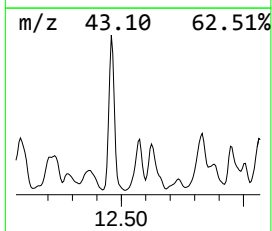
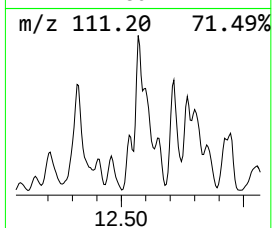
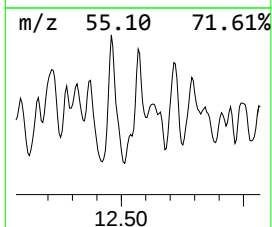
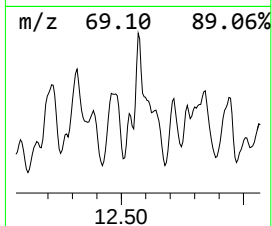
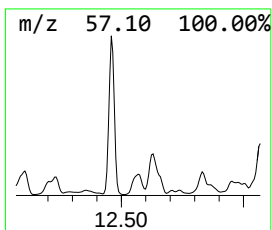
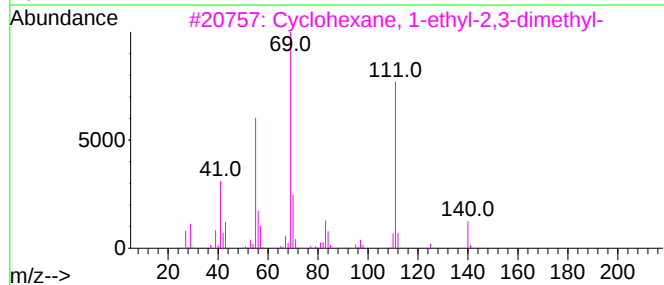
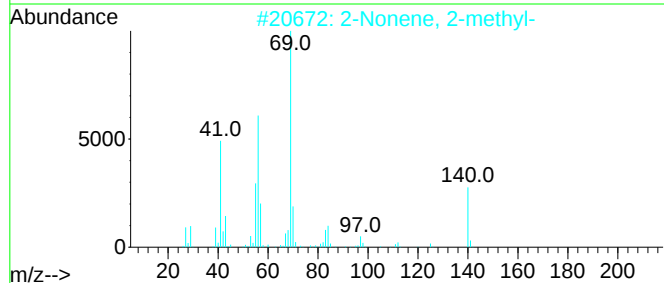
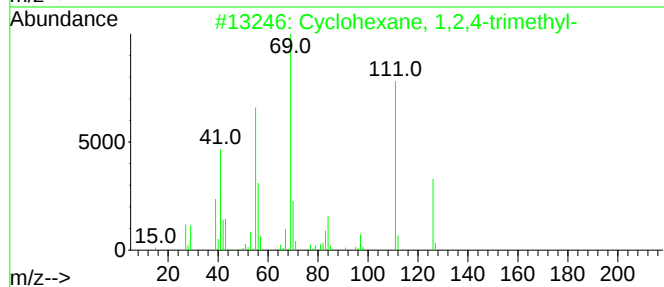
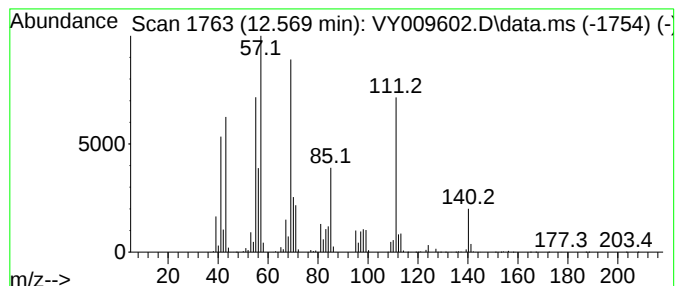
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown12.569 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.569	128.62 ug/l	5167990	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	47
2		2-Nonene, 2-methyl-	140	C10H20	002129-95-5	46
3		Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	46
4		Ether, heptyl hexyl	200	C13H28O	007289-40-9	38
5		Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

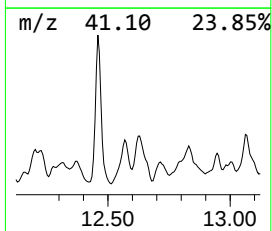
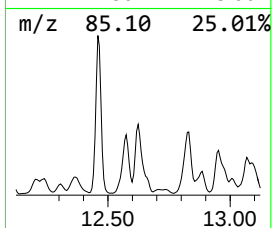
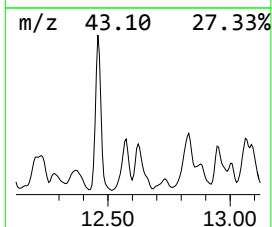
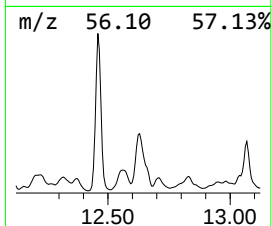
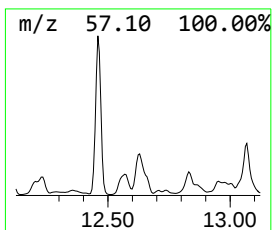
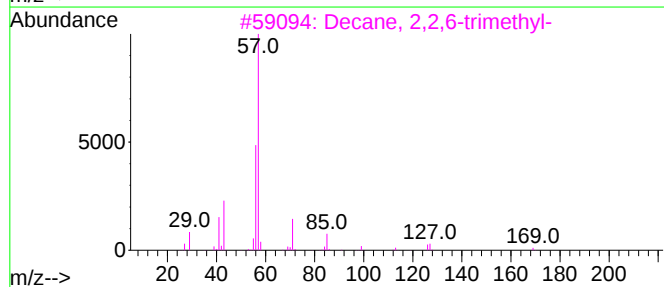
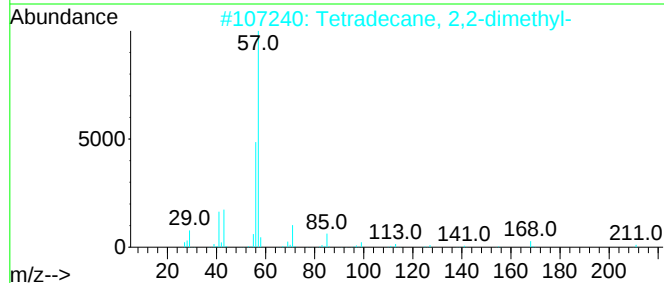
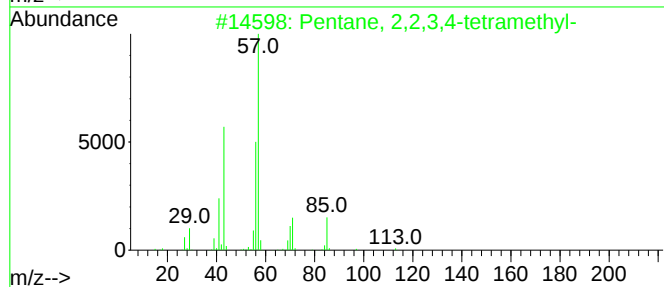
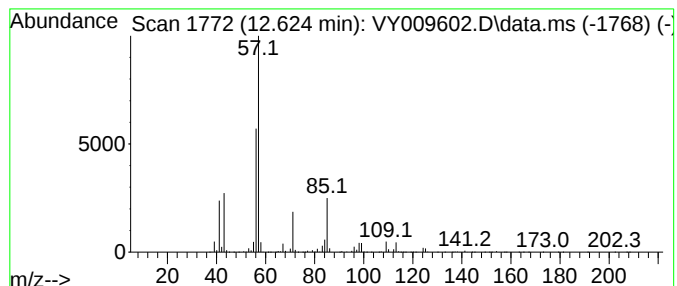
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Pentane, 2,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.624	134.59 ug/l	5407710	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	59
2			Tetradecane, 2,2-dimethyl-	226	C16H34	059222-86-5	59
3			Decane, 2,2,6-trimethyl-	184	C13H28	062237-97-2	53
4			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	53
5			Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

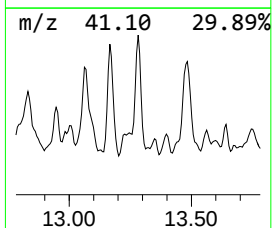
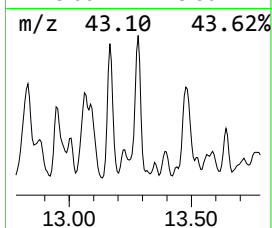
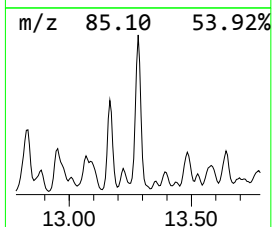
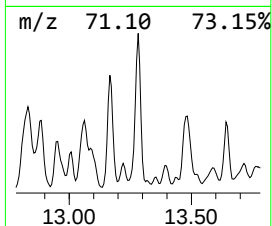
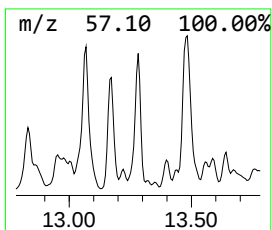
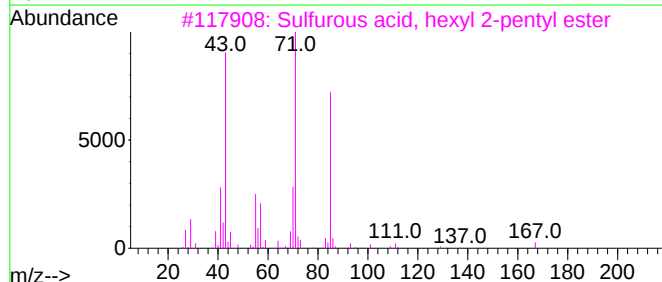
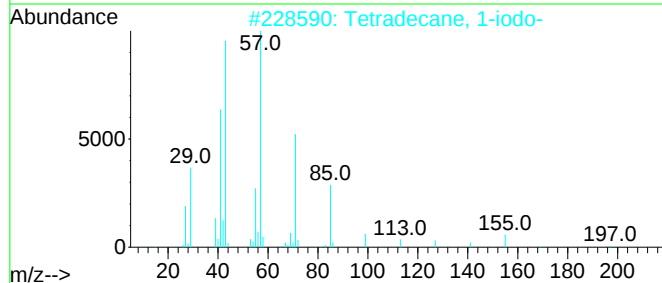
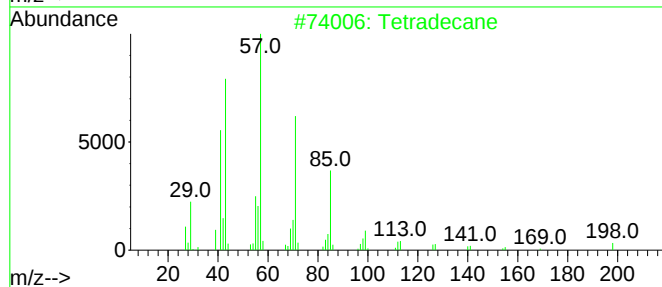
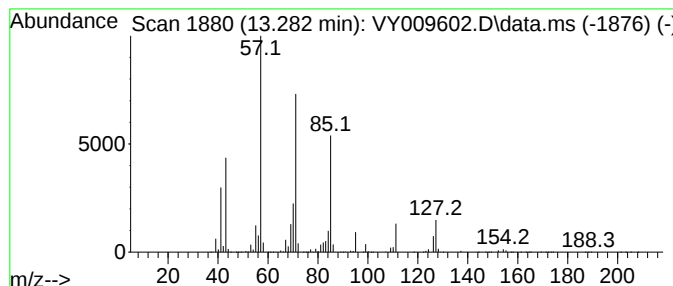
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.282	110.01 ug/l	4420060	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	59
2			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	53
3			Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	53
4			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	53
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

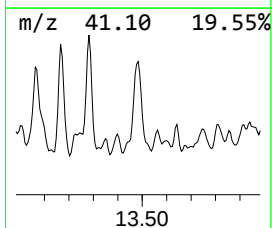
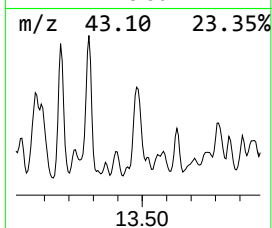
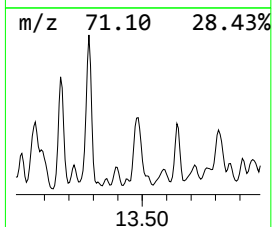
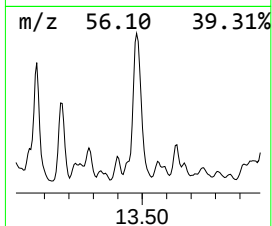
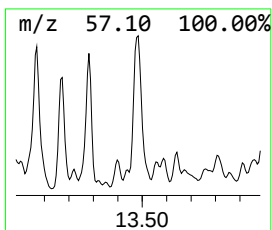
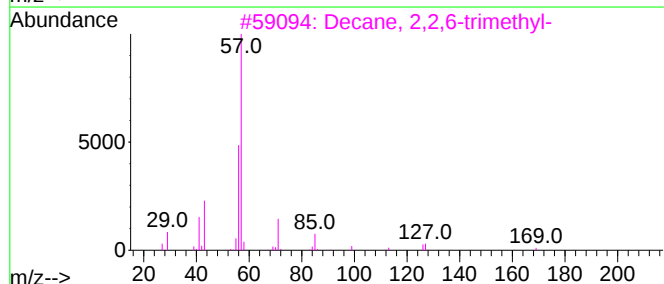
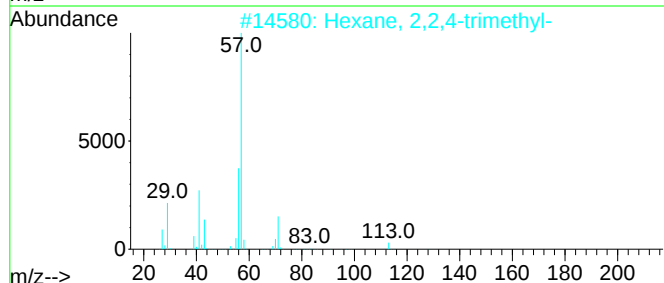
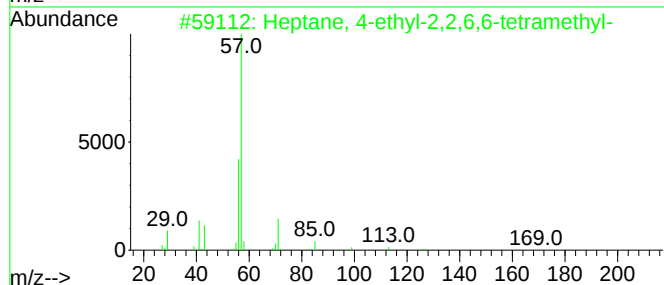
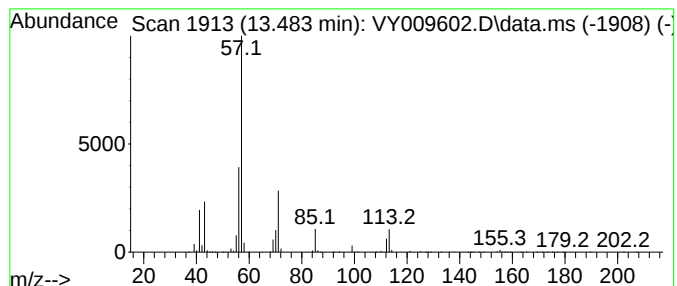
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Heptane, 4-ethyl-2,2,6,6-te... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.483	103.02 ug/l	4139500	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-ethyl-2,2,6,6-tetrame...	184	C13H28	062108-31-0	72
2			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
3			Decane, 2,2,6-trimethyl-	184	C13H28	062237-97-2	53
4			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53
5			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

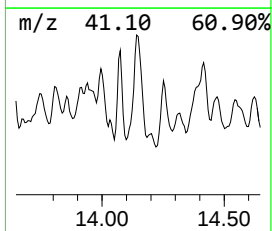
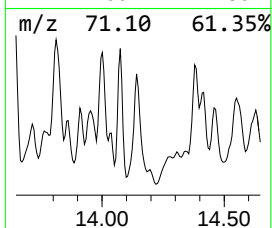
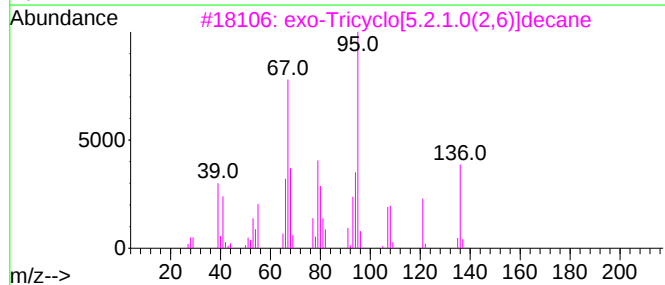
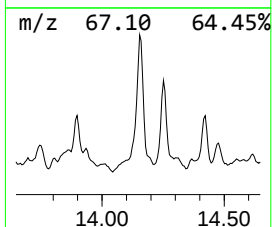
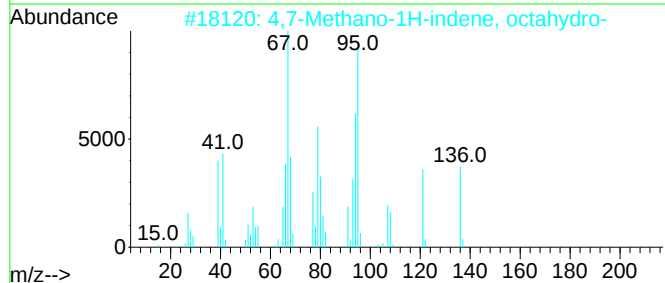
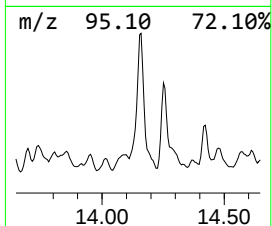
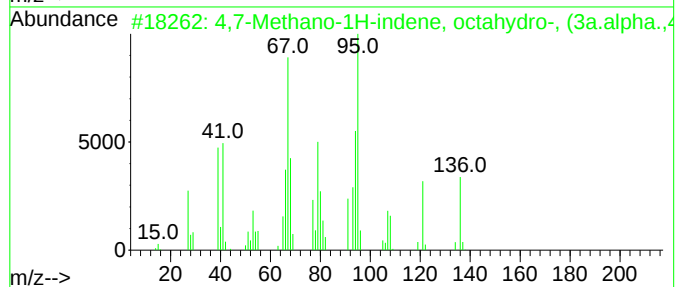
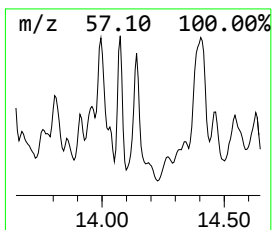
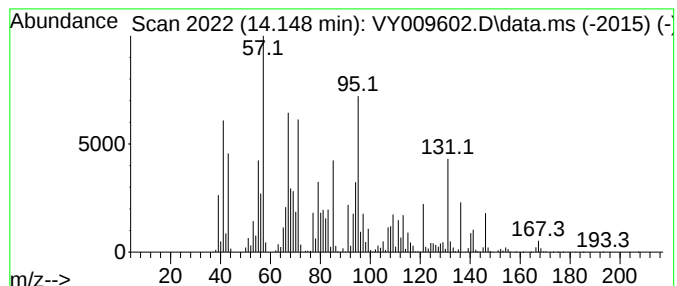
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 4,7-Methano-1H-indene, octa... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.148	121.34 ug/l	4875370	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,7-Methano-1H-indene, octahydro...	136	C10H16	002825-83-4	96
2			4,7-Methano-1H-indene, octahydro-	136	C10H16	006004-38-2	96
3			exo-Tricyclo[5.2.1.0(2,6)]decane	136	C10H16	002825-82-3	87
4			1-Cyclopentylcyclopentene	136	C10H16	004884-21-3	50
5			Undecane, 3-methyl-	170	C12H26	001002-43-3	15



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY071922\
Data File : VY009602.D
Acq On : 19 Jul 2022 20:10
Operator : KP/MD
Sample : N3652-05
Misc : 7.39g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SED-3-(6-12)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071822S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 2,2,4-...	8.319	2792.8	ug/l	60093600	2	8.685	1075880	50.0
Pentane, 2,2,3-...	9.319	286.4	ug/l	6162010	2	8.685	1075880	50.0
Pentane, 2,3,4-...	9.618	2392.6	ug/l	51481900	2	8.685	1075880	50.0
Pentane, 2,3,3-...	9.746	2844.1	ug/l	61198500	2	8.685	1075880	50.0
Hexane, 2,2,5-t...	10.112	226.6	ug/l	20694500	3	11.490	4566230	50.0
unknown12.569	12.569	128.6	ug/l	5167990	4	13.422	2008990	50.0
Pentane, 2,2,3,...	12.624	134.6	ug/l	5407710	4	13.422	2008990	50.0
Tetradecane	13.282	110.0	ug/l	4420060	4	13.422	2008990	50.0
Heptane, 4-ethy...	13.483	103.0	ug/l	4139500	4	13.422	2008990	50.0
4,7-Methano-1H-...	14.148	121.3	ug/l	4875370	4	13.422	2008990	50.0