

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
 Data File : VY017223.D
 Acq On : 05 Feb 2024 12:56
 Operator : SY/MD
 Sample : P1366-03
 Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 EB-1-0-2

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\82Y020124S.M

Title : SW846 8260

Signal : TIC: VY017223.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.704	786	801	815	rBV	278463	872324	6.54%	1.549%
2	7.728	960	969	974	rBV2	61886	160435	1.20%	0.285%
3	7.795	974	980	986	rVV2	138700	361724	2.71%	0.642%
4	7.886	986	995	1008	rVV	962727	2414050	18.11%	4.287%
5	8.338	1051	1069	1082	rBV	5893539	13328908	100.00%	23.668%
6	8.703	1120	1129	1141	rBV	222350	447753	3.36%	0.795%
7	9.203	1198	1211	1215	rBV	729344	1448203	10.87%	2.572%
8	9.252	1215	1219	1226	rVV	877793	1645189	12.34%	2.921%
9	9.331	1226	1232	1246	rVB	511352	1038642	7.79%	1.844%
10	9.630	1271	1281	1292	rBV	4074252	7382060	55.38%	13.108%
11	9.764	1292	1303	1310	rBV	2994699	6211621	46.60%	11.030%
12	9.953	1326	1334	1349	rVB2	166395	377900	2.84%	0.671%
13	10.124	1350	1362	1368	rBV	2122455	3578126	26.84%	6.354%
14	10.191	1368	1373	1378	rBV	262086	413776	3.10%	0.735%
15	10.258	1378	1384	1389	rBV	800123	1360051	10.20%	2.415%
16	10.648	1441	1448	1455	rVB	390181	657270	4.93%	1.167%
17	10.916	1485	1492	1500	rBV	95708	185832	1.39%	0.330%
18	11.495	1579	1587	1593	rBV2	556249	1299207	9.75%	2.307%
19	11.575	1593	1600	1616	rVB2	484885	982614	7.37%	1.745%
20	11.715	1616	1623	1625	rBV	78998	144898	1.09%	0.257%
21	11.758	1625	1630	1641	rVB	330945	551229	4.14%	0.979%
22	11.947	1657	1661	1665	rBV	155039	244044	1.83%	0.433%
23	12.123	1684	1690	1697	rVB	182071	300697	2.26%	0.534%
24	12.471	1741	1747	1757	rBV	1886479	3044549	22.84%	5.406%
25	12.568	1757	1763	1769	rBV2	116342	271473	2.04%	0.482%
26	12.636	1769	1774	1785	rVB	427961	878279	6.59%	1.560%
27	12.745	1785	1792	1795	rBV2	81874	135686	1.02%	0.241%
28	12.849	1800	1809	1814	rBV	489108	1001258	7.51%	1.778%
29	12.898	1814	1817	1825	rVB	82690	135465	1.02%	0.241%
30	13.081	1839	1847	1852	rVB	520744	850755	6.38%	1.511%
31	13.190	1860	1865	1870	rVB	176182	267845	2.01%	0.476%
32	13.294	1870	1882	1886	rBV	453248	964612	7.24%	1.713%
33	13.440	1897	1906	1910	rBV2	369063	793011	5.95%	1.408%
34	13.489	1910	1914	1924	rVB	606147	1327254	9.96%	2.357%
35	13.654	1934	1941	1949	rBV3	247497	459436	3.45%	0.816%
36	13.837	1967	1971	1977	rVB2	78513	140950	1.06%	0.250%

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ClientSampleId :
EB-1-0-2

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\82Y020124S.M
Title : SW846 8260

37	14.001	1988	1998	2008	rVB2	96359	315368	2.37%	0.560%
38	14.397	2057	2063	2073	rVB3	149898	324266	2.43%	0.576%

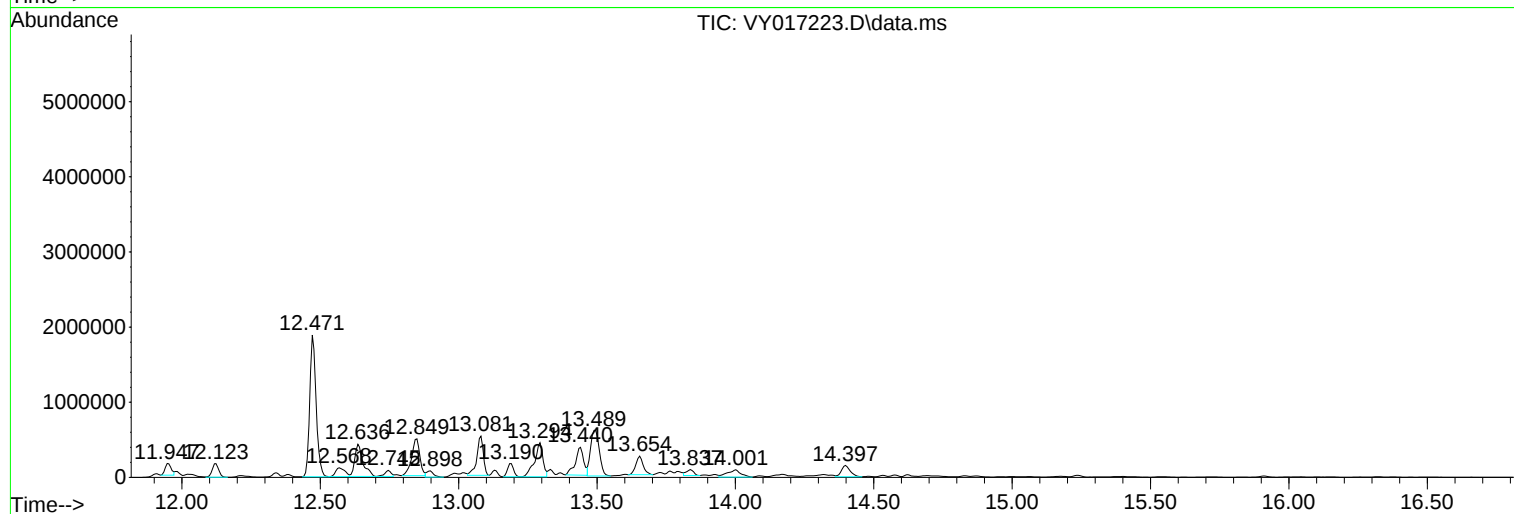
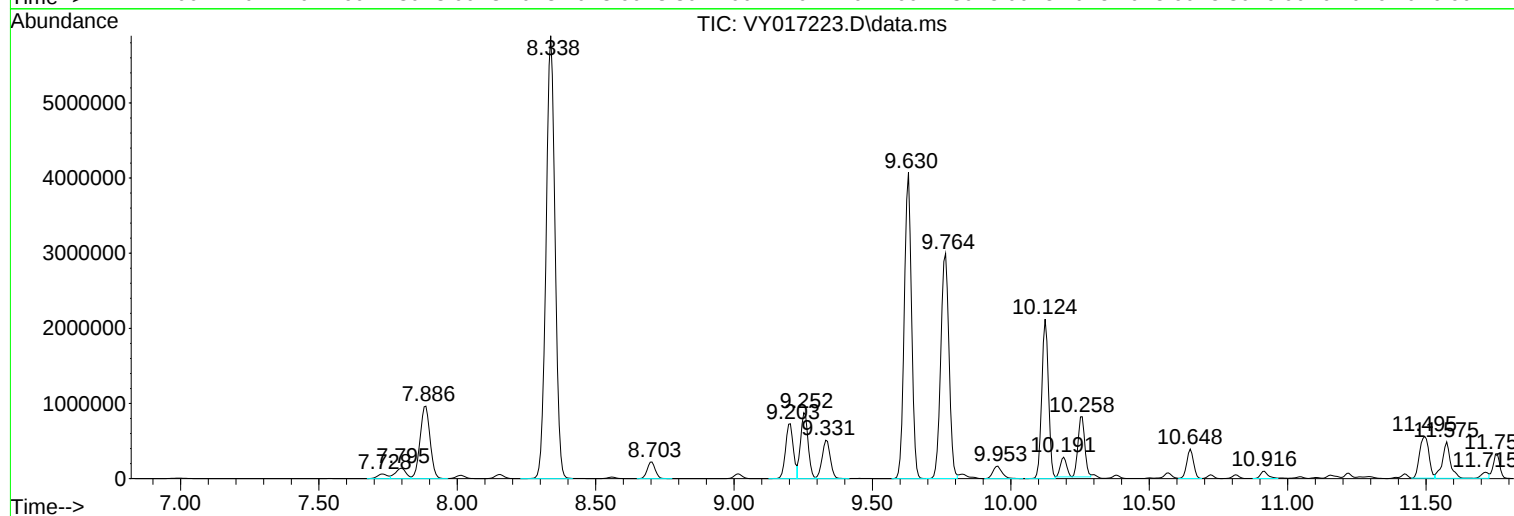
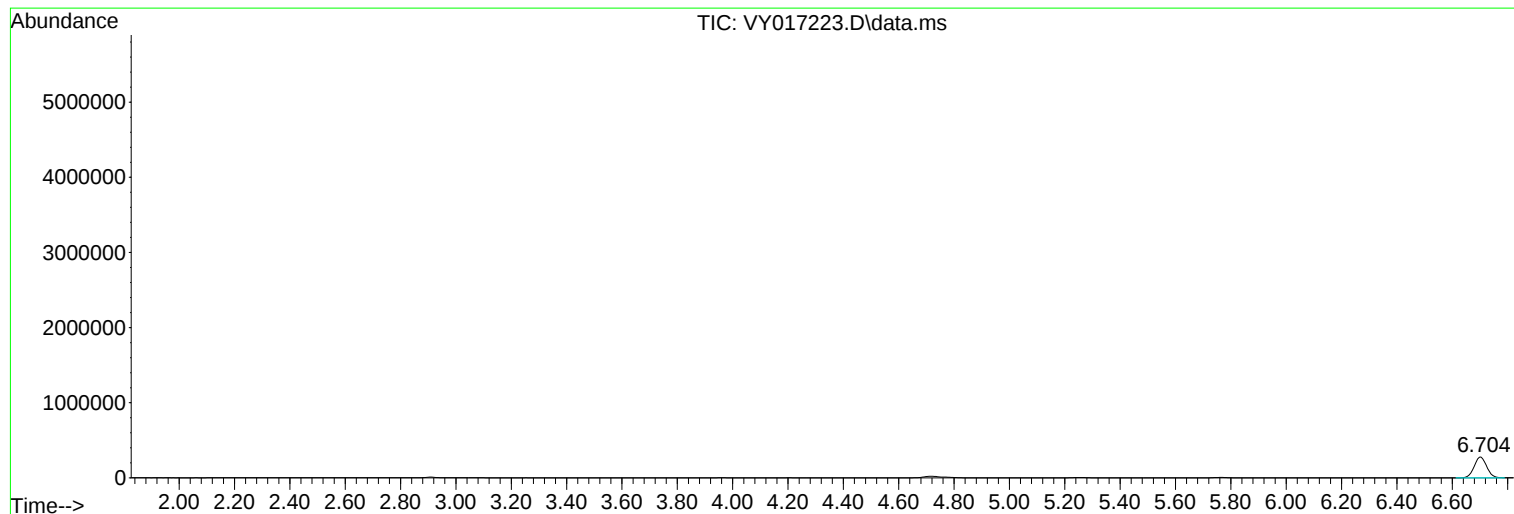
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

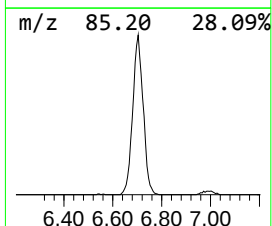
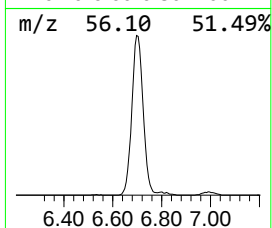
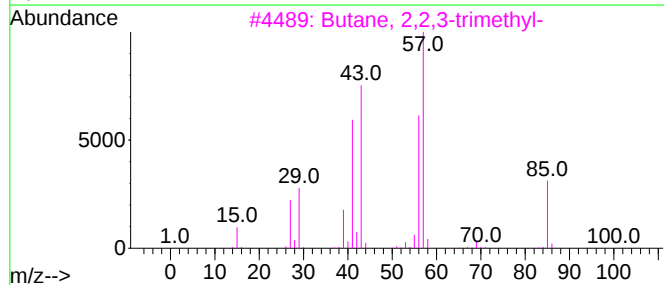
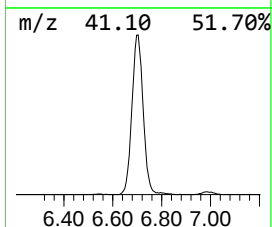
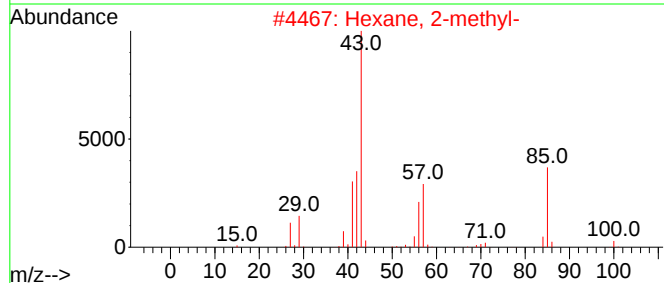
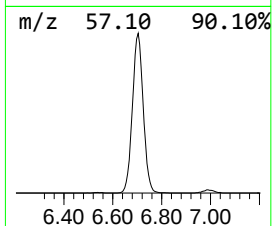
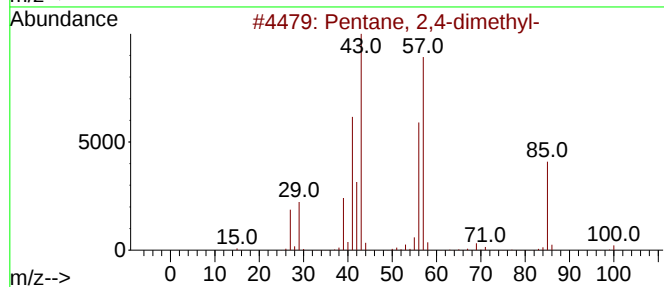
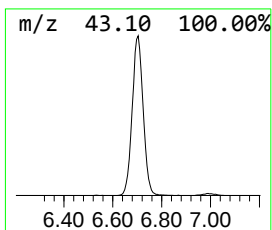
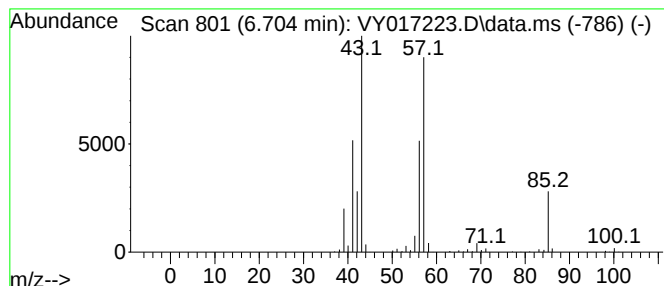
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.704	120.58 ug/l	872324	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	94
2			Hexane, 2-methyl-	100	C7H16	000591-76-4	72
3			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59
4			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	59
5			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	50



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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

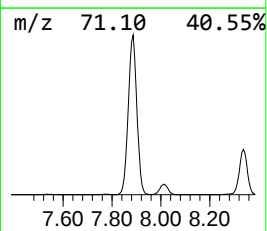
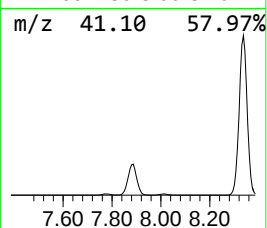
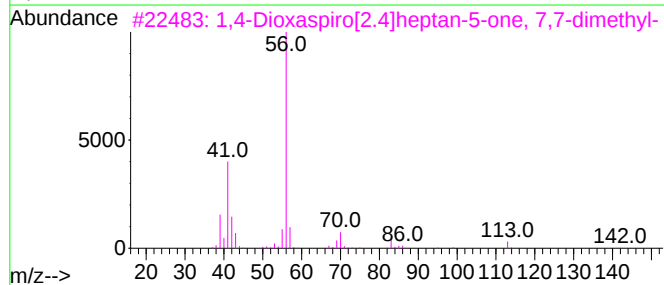
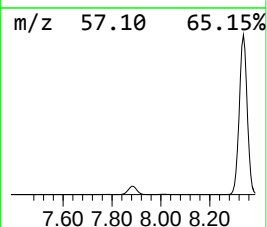
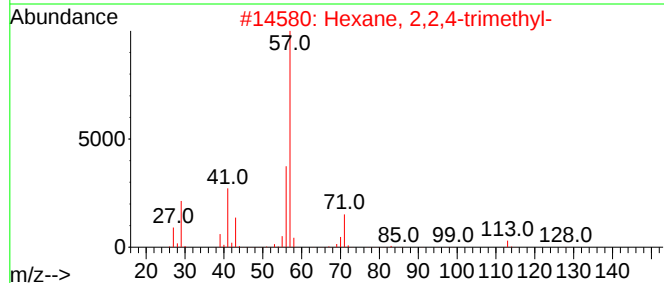
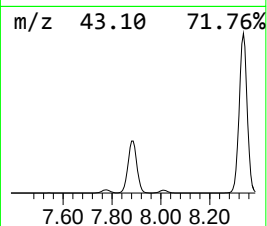
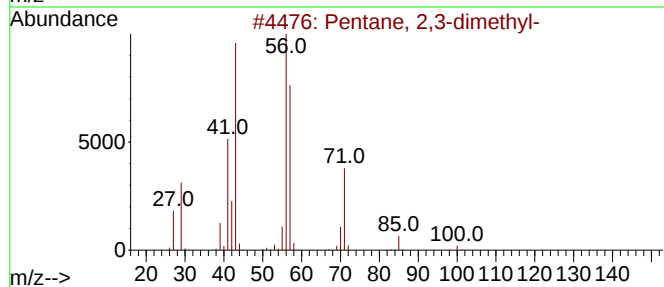
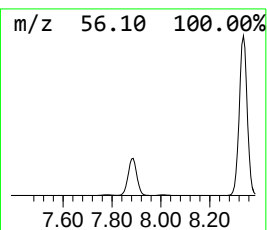
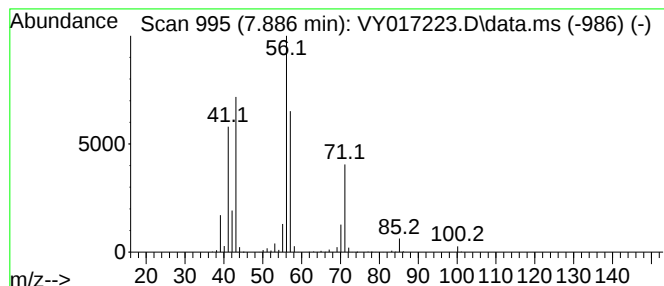
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.886	333.69 ug/l	2414050	Pentafluorobenzene	7.801

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45
3			1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	40
4			Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	38
5			Aziridine, 2-methyl-	57	C3H7N	000075-55-8	38



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Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

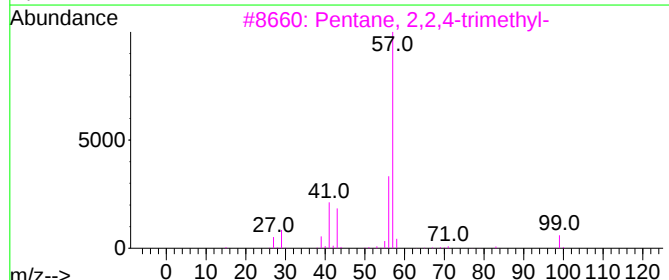
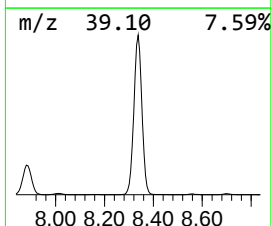
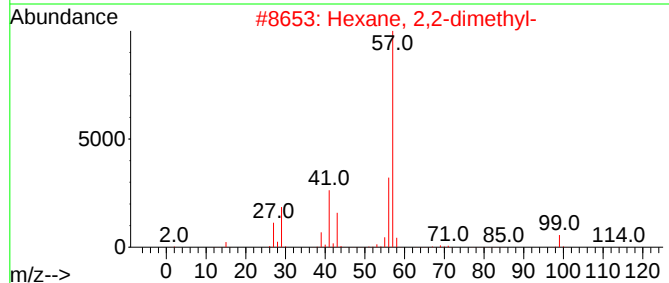
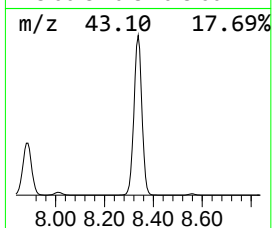
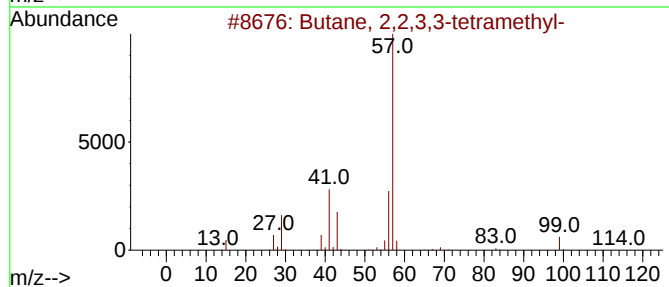
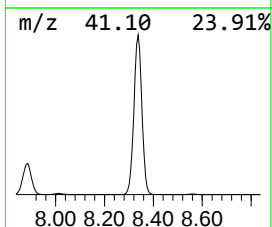
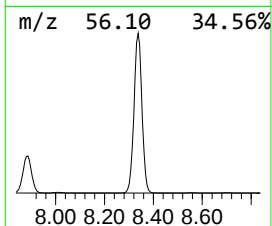
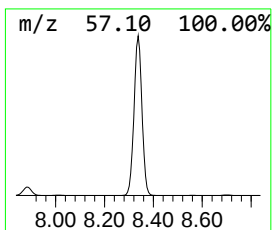
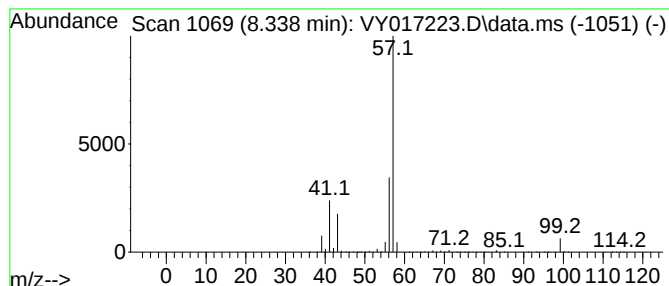
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 3 Butane, 2,2,3,3-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.338	1488.42 ug/l	13328900	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	78
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64



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Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

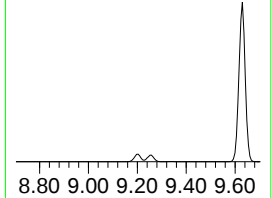
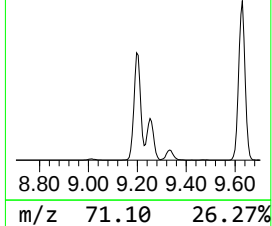
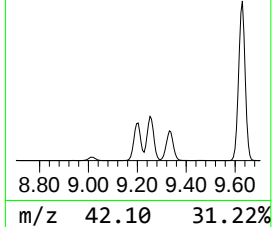
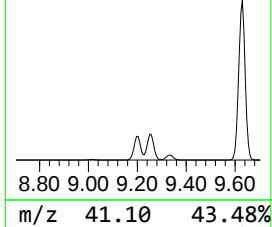
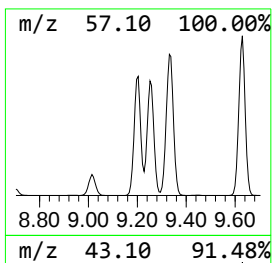
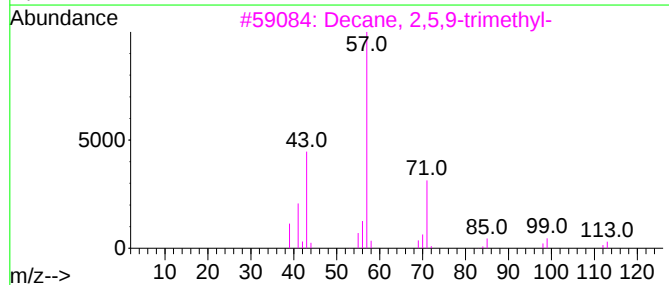
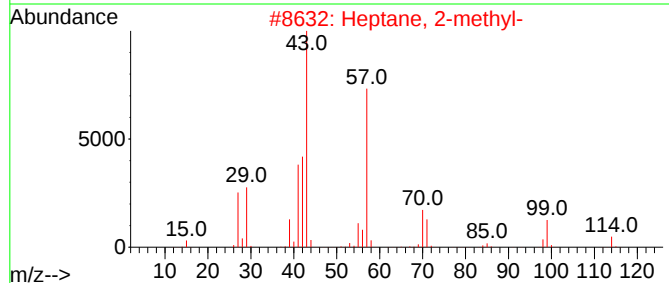
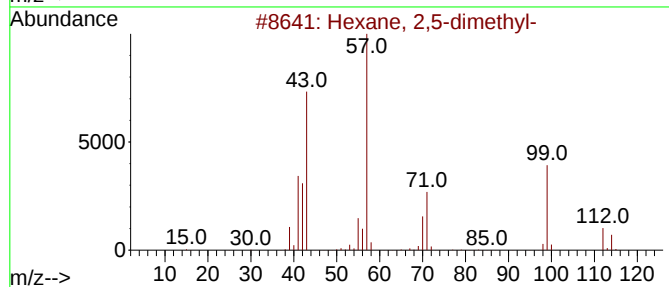
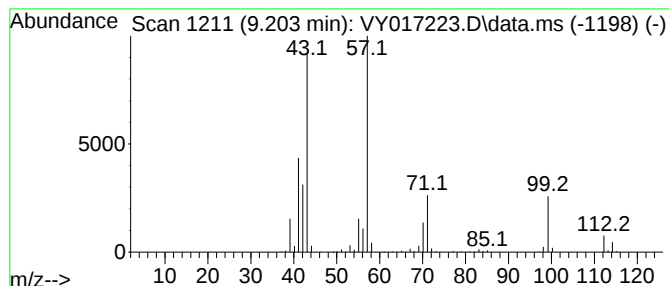
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 4 Hexane, 2,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.203	161.72 ug/l	1448200	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	94
2			Heptane, 2-methyl-	114	C8H18	000592-27-8	52
3			Decane, 2,5,9-trimethyl-	184	C13H28	062108-22-9	50
4			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
5			(Aminomethyl)cyclopropane	71	C4H9N	002516-47-4	47



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Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

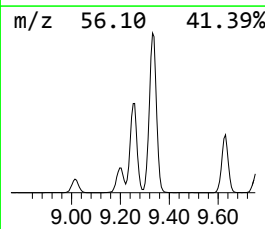
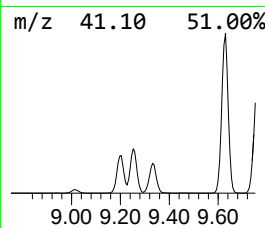
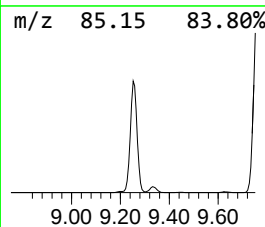
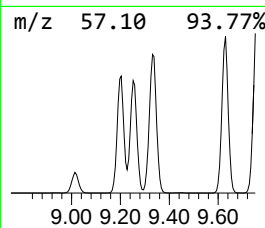
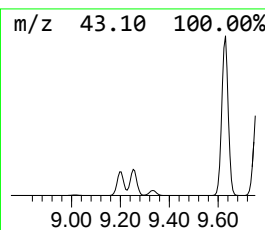
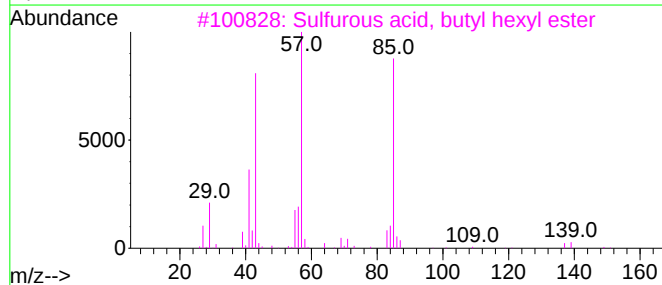
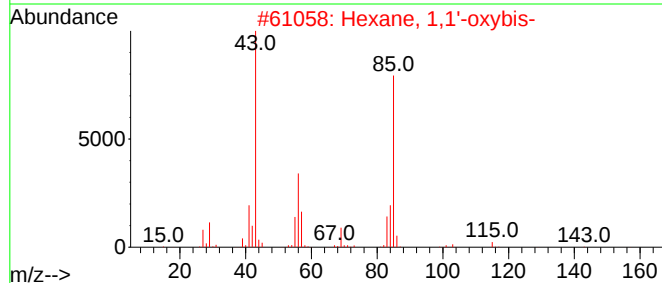
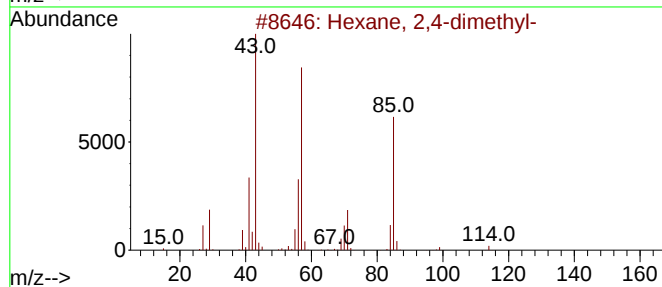
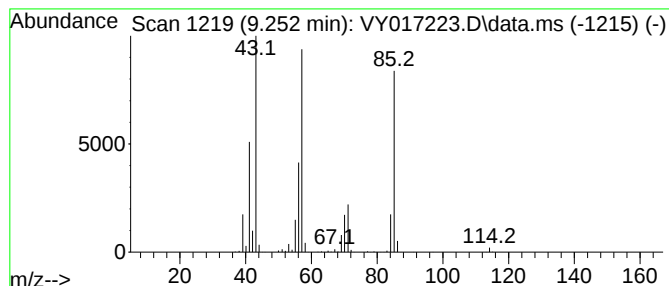
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Quant Title : SW846 8260

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

Peak Number 5 Hexane, 2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.252	183.72 ug/l	1645190	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
2			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	64
3			Sulfurous acid, butyl hexyl ester	222	C10H22O3S	1000309-17-3	59
4			Sulfurous acid, butyl isohexyl e...	222	C10H22O3S	1000309-17-2	53
5			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

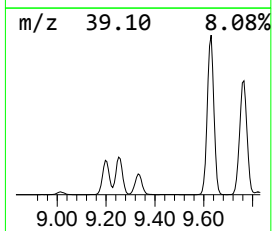
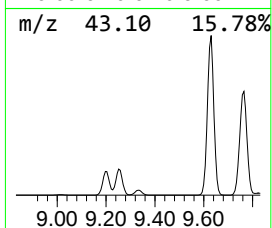
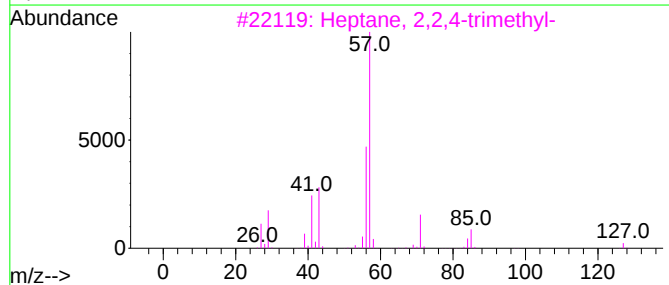
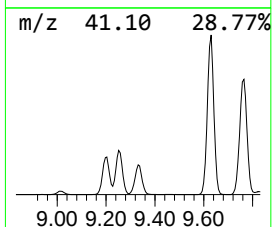
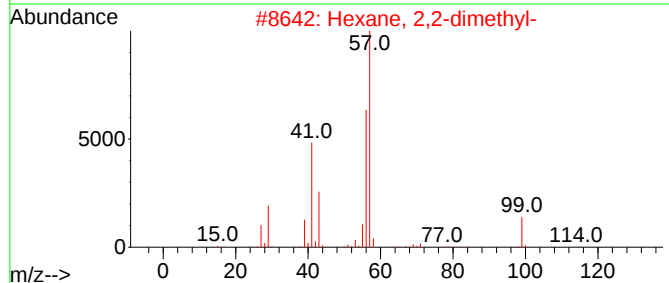
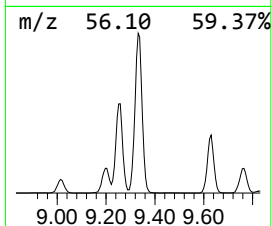
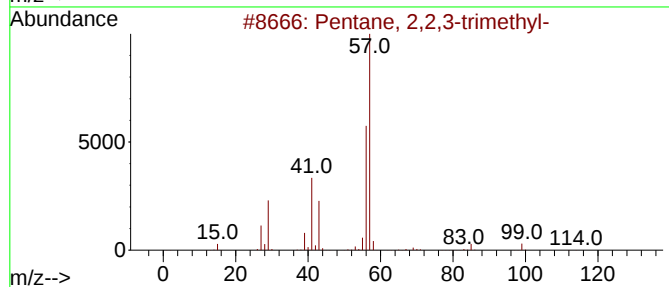
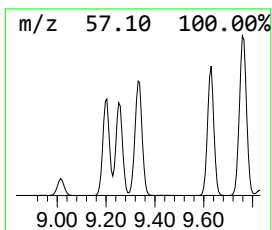
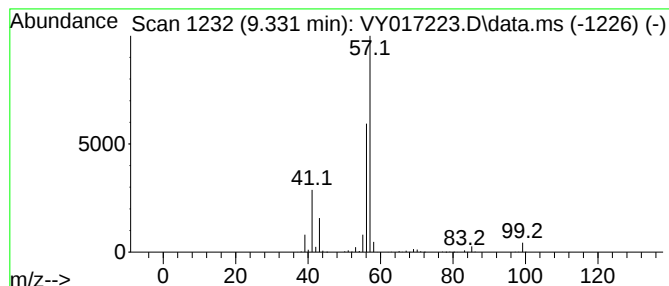
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Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.331	115.98 ug/l	1038640	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	83
2			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	78
4			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5			Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

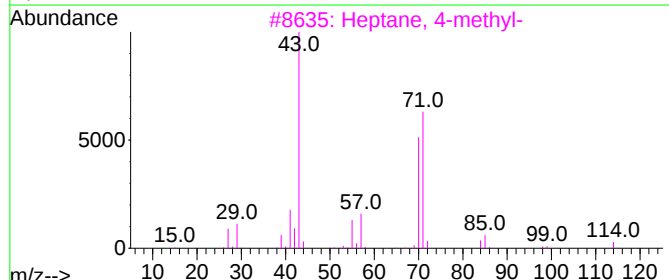
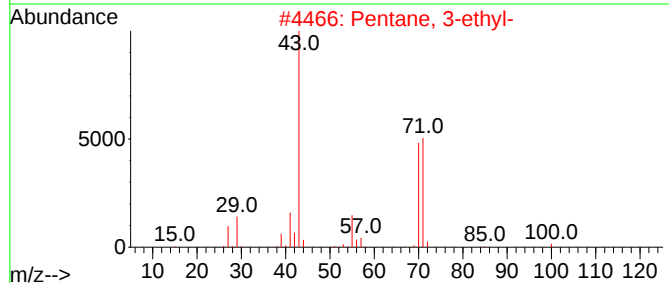
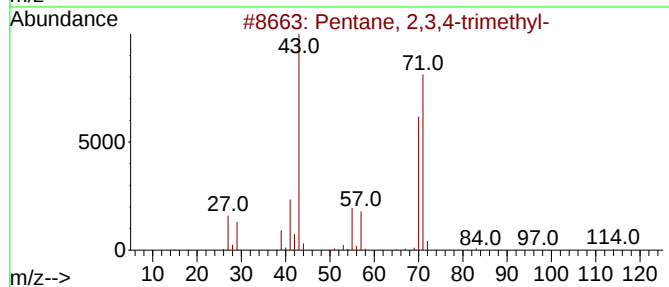
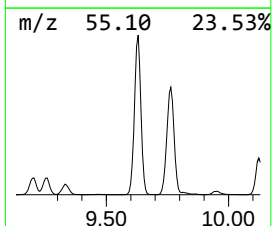
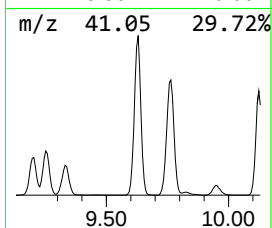
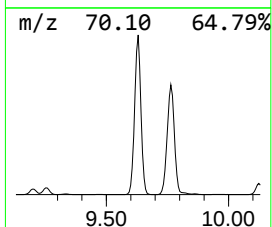
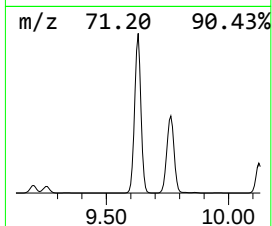
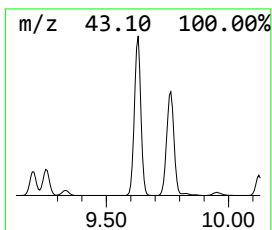
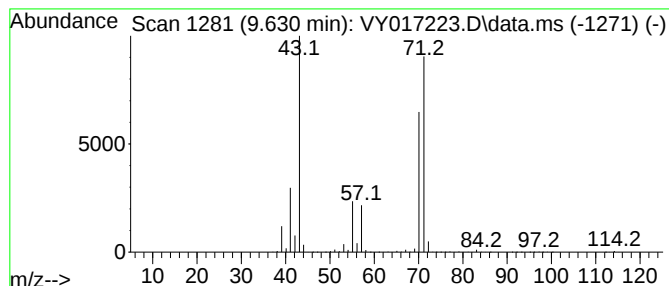
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.630	824.35 ug/l	7382060	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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	72
5			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

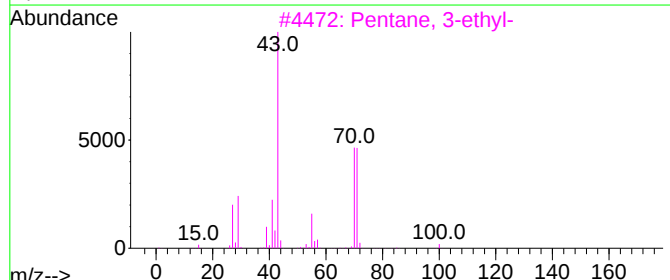
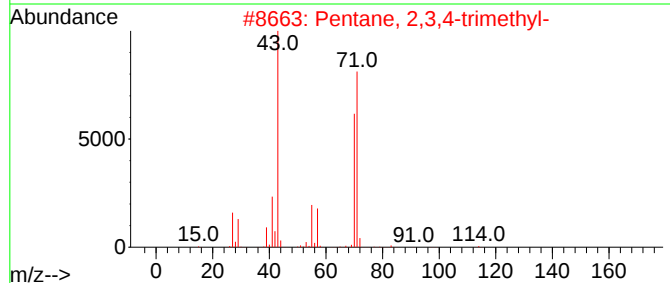
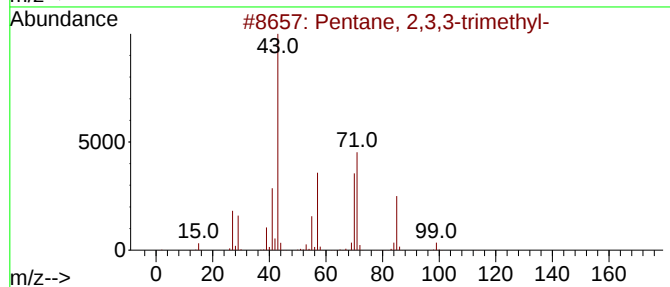
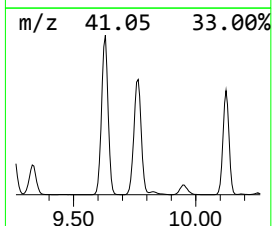
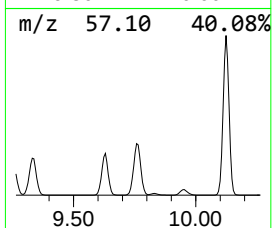
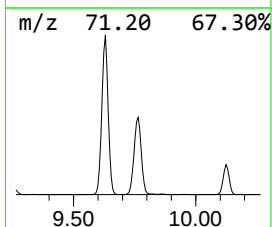
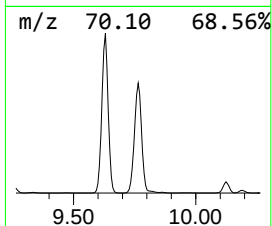
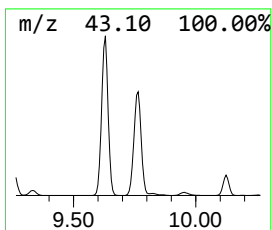
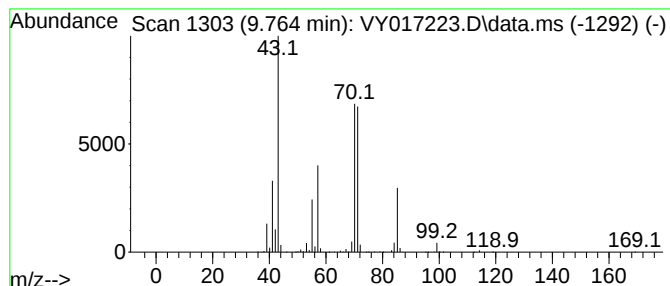
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.764	693.64 ug/l	6211620	1,4-Difluorobenzene	8.703

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
2			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
3			Pentane, 3-ethyl-	100	C7H16	000617-78-7	59
4			Diisoamyl ether	158	C10H22O	000544-01-4	59
5			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

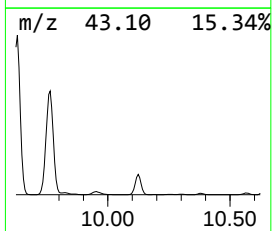
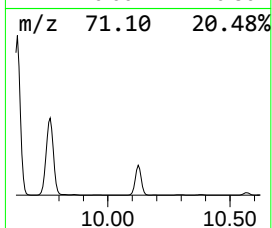
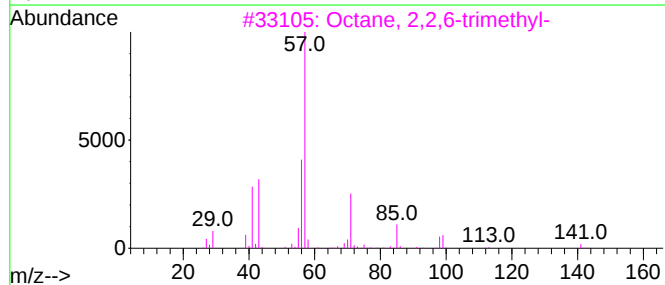
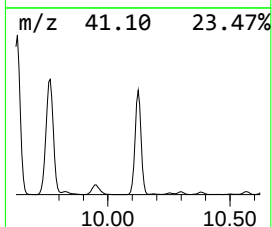
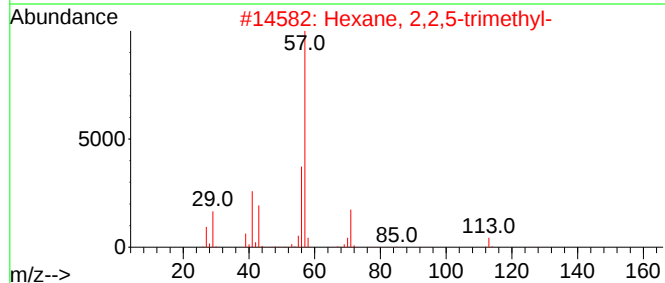
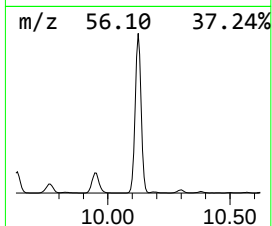
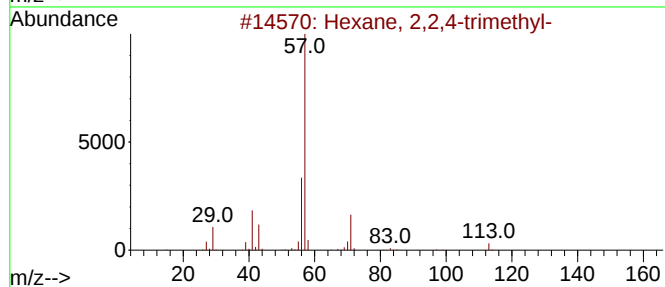
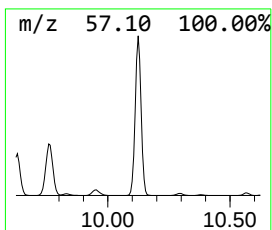
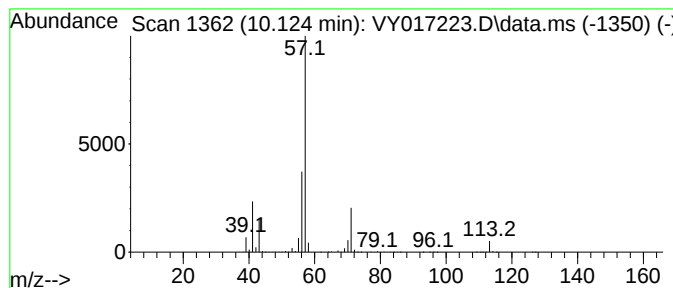
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Quant Title : SW846 8260

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

Peak Number 9 Hexane, 2,2,4-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.124	137.70 ug/l	3578130	Chlorobenzene-d5	11.502

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	83
2			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	83
3			Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	78
4			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

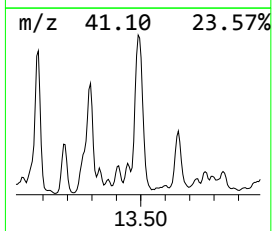
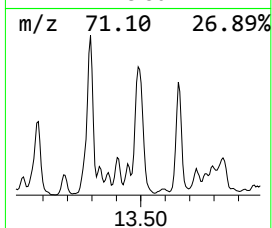
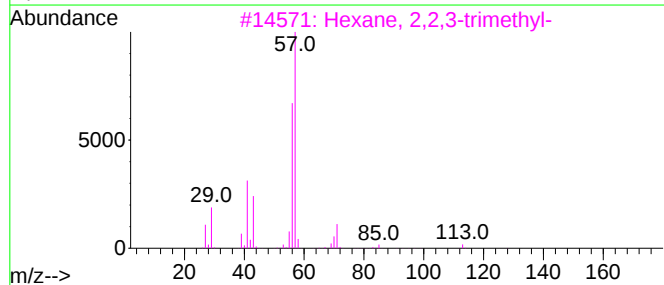
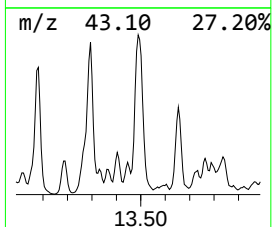
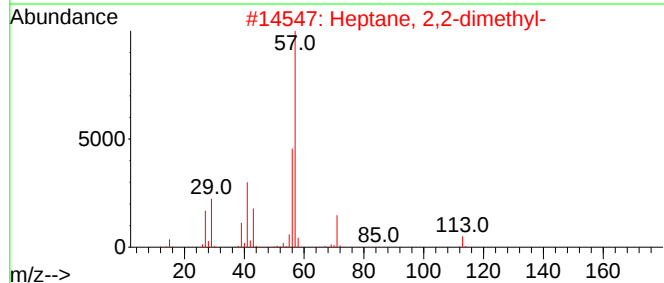
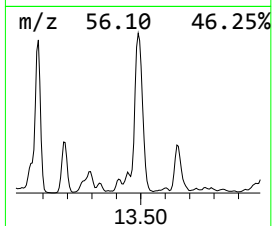
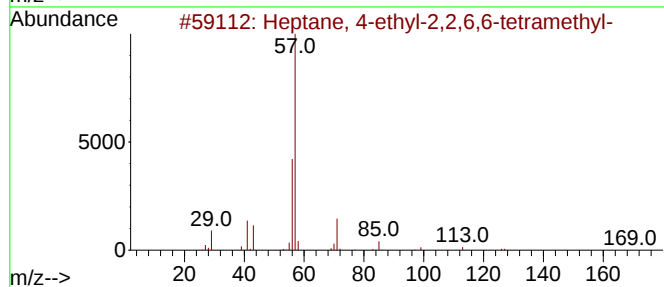
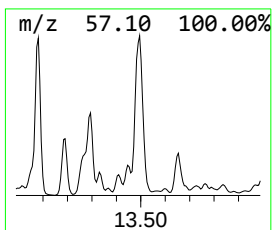
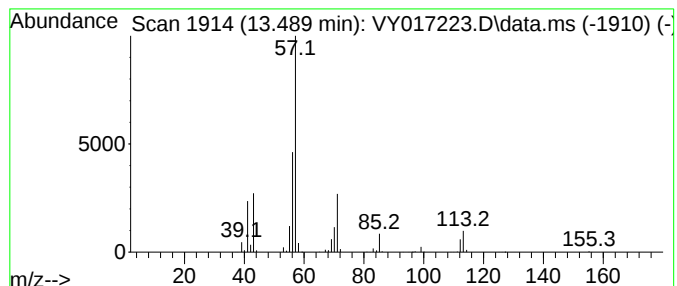
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.489	83.68 ug/l	1327250	1,4-Dichlorobenzene-d4	13.434

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			Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64
3			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4			Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	64
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020524\
Data File : VY017223.D
Acq On : 05 Feb 2024 12:56
Operator : SY/MD
Sample : P1366-03
Misc : 5.87g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EB-1-0-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.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, 2,4-di...	6.704	120.6	ug/l	872324	1	7.801	361724	50.0
Pentane, 2,3-di...	7.886	333.7	ug/l	2414050	1	7.801	361724	50.0
Butane, 2,2,3,3...	8.338	1488.4	ug/l	13328900	2	8.703	447753	50.0
Hexane, 2,5-dim...	9.203	161.7	ug/l	1448200	2	8.703	447753	50.0
Hexane, 2,4-dim...	9.252	183.7	ug/l	1645190	2	8.703	447753	50.0
Pentane, 2,2,3-...	9.331	116.0	ug/l	1038640	2	8.703	447753	50.0
Pentane, 2,3,4-...	9.630	824.4	ug/l	7382060	2	8.703	447753	50.0
Pentane, 2,3,3-...	9.764	693.6	ug/l	6211620	2	8.703	447753	50.0
Hexane, 2,2,4-t...	10.124	137.7	ug/l	3578130	3	11.502	1299210	50.0
Heptane, 4-ethy...	13.489	83.7	ug/l	1327250	4	13.434	793011	50.0