LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

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\82Y062325S.M
Title : SW846 8260

Signal : TIC: VY022899.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.616	463	475	487	rBV3	29325	94562	4.93%	0.466%
2	7.634	960	970	976	rBV	230313	557565	29.08%	2.745%
3	7.707	976	982	991	rVV	309832	709417	37.01%	3.493%
4	8.061	1032	1040	1053	rBV	229013	506572	26.42%	2.494%
5	8.615	1122	1131	1141	rBV	575607	1133631	59.13%	5.582%
6	9.109	1198	1212	1217	rBV3	44771	111569	5.82%	0.549%
7	9.170	1217	1222	1229	rVB2	32814	61923	3.23%	0.305%
8	9.359	1246	1253	1267	rBV2	65160	164076	8.56%	0.808%
9	9.536	1273	1282	1291	rBV4	43352	95743	4.99%	0.471%
10	9.688	1294	1307	1311	rBV	142908	310702	16.21%	1.530%
11	9.731	1311	1314	1320	rVB2	76477	128595	6.71%	0.633%
12	9.865	1327	1336	1340	rBV	250876	490761	25.60%	2.416%
13	9.920	1340	1345	1353	rVB	282097	545420	28.45%	2.686%
14	10.030	1353	1363	1368	rBV5	41985	120886	6.31%	0.595%
15	10.103	1368	1375	1389	rVB	1048986	1917042	100.00%	9.439%
16	10.231	1389	1396	1398	rBV4	22477	40400	2.11%	0.199%
17	10.274	1398	1403	1405	rBV2	92311	163469	8.53%	0.805%
18	10.377	1413	1420	1423	rBV	45651	84518	4.41%	0.416%
19	10.444	1427	1431	1440	rVV	94973	182235	9.51%	0.897%
20	10.548	1440	1448	1453	rVV5	97302	271973	14.19%	1.339%
21	10.591	1453	1455	1459	rVV2	62466	90937	4.74%	0.448%
22	10.639	1459	1463	1468	rVV	105346	168123	8.77%	0.828%
23	10.731	1472	1478	1483	rVV2	113640	225608	11.77%	1.111%
24	10.780	1483	1486	1489	rVV2	30261	50616	2.64%	0.249%
25	10.835	1489	1495	1502	rVV	162001	385300	20.10%	1.897%
26	10.902	1502	1506	1509	rVV	61207	114929	6.00%	0.566%
27	10.963	1513	1516	1520	rVV2	48656	79364	4.14%	0.391%
28	11.017	1520	1525	1530	rVV2	140174	258155	13.47%	1.271%
29	11.072	1530	1534	1538	rVV2	37717	70122	3.66%	0.345%
30	11.133	1538	1544	1547	rVV4	72514	149627	7.81%	0.737%
31	11.170	1547	1550	1555	rVV	131887	227514	11.87%	1.120%
32	11.219	1555	1558	1566	rVB4	33490	73523	3.84%	0.362%
33	11.304	1566	1572	1577	rBV	52422	87744	4.58%	0.432%
34	11.414	1583	1590	1600	rBV	951390	1551422	80.93%	7.639%
35	11.505	1600	1605	1609	rVV4	35390	67145	3.50%	0.331%
36	11.554	1609	1613	1616	rVV5	16929	26312	1.37%	0.130%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
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Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

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\82Y062325S.M
Title : SW846 8260

37	11.603	1616	1621	1625	rVV7	17438	43000	2.24%	0.212%
38	11.664	1625	1631	1642	rVB5	48761	144909	7.56%	0.713%
39	11.822	1651	1657	1661	rBV2	13191	22287	1.16%	0.110%
40	11.926	1666	1674	1677	rBV6	37336	92374	4.82%	0.455%
41	12.127	1702	1707	1710	rVV3	33351	56443	2.94%	0.278%
42	12.164	1710	1713	1717	rVB3	15757	28696	1.50%	0.141%
43	12.407	1747	1753	1759	rBV	707574	1113799	58.10%	5.484%
44	12.481	1760	1765	1771	rBV5	16554	36887	1.92%	0.182%
45	12.566	1771	1779	1786	rVB2	89253	171369	8.94%	0.844%
46	12.651	1786	1793	1798	rBV6	27188	72830	3.80%	0.359%
47	12.731	1799	1806	1811	rBV4	66323	131621	6.87%	0.648%
48	12.871	1825	1829	1836	rBV3	31884	83696	4.37%	0.412%
49	12.962	1840	1844	1852	rVV5	67026	121323	6.33%	0.597%
50	13.048	1853	1858	1861	rVV7	22303	35565	1.86%	0.175%
51	13.090	1861	1865	1871	rVV6	18089	44028	2.30%	0.217%
52	13.169	1875	1878	1882	rVB3	29308	40024	2.09%	0.197%
53	13.237	1882	1889	1893	rBV5	49832	104332	5.44%	0.514%
54	13.285	1893	1897	1901	rVB5	17529	30471	1.59%	0.150%
55	13.346	1901	1907	1916	rVB	760279	1250492	65.23%	6.157%
56	13.450	1920	1924	1925	rBV	17214	21445	1.12%	0.106%
57	13.486	1926	1930	1936	rBV4	56322	86598	4.52%	0.426%
58	13.584	1942	1946	1953	rBV2	64997	110461	5.76%	0.544%
59	13.663	1953	1959	1965	rBV3	187758	348238	18.17%	1.715%
60	13.718	1965	1968	1969	rBV2	16545	19275	1.01%	0.095%
61	13.743	1969	1972	1975	rBV	30136	31473	1.64%	0.155%
62	13.797	1978	1981	1984	rBV3	35564	55425	2.89%	0.273%
63	13.834	1984	1987	1992	rVB6	54859	75851	3.96%	0.373%
64	13.889	1992	1996	2000	rVB2	119155	170777	8.91%	0.841%
65	13.932	2001	2003	2008	rVB5	35997	46020	2.40%	0.227%
66	13.992	2008	2013	2018	rBV2	227831	394176	20.56%	1.941%
67	14.047	2018	2022	2030	rVB4	102747	237265	12.38%	1.168%
68	14.133	2030	2036	2039	rBV	118174	205363	10.71%	1.011%
69	14.181	2039	2044	2047	rBV3	113543	237744	12.40%	1.171%
70	14.255	2052	2056	2059	rVB	112081	172455	9.00%	0.849%
71	14.297	2059	2063	2066	rBV2	124907	177302	9.25%	0.873%
72	14.340	2066	2070	2073	rBV2	127166	164723	8.59%	0.811%
73	14.486	2090	2094	2098	rBV5	37648	71195	3.71%	0.351%
74	14.529	2098	2101	2106	rBV3	100381	141167	7.36%	0.695%
75	14.602	2106	2113	2117	rBV5	141455	356504	18.60%	1.755%
76	14.651	2117	2121	2127	rVB5	113430	188908	9.85%	0.930%

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ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

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
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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Title : SW846 8260

77	14.706	2127	2130	2134	rVB2	49182	71086	3.71%	0.350%
78	14.748	2135	2137	2140	rBV2	29800	33392	1.74%	0.164%
79	14.785	2140	2143	2147	rVB4	53933	71354	3.72%	0.351%
80	14.822	2147	2149	2151	rBV3	27278	28665	1.50%	0.141%
81	14.858	2151	2155	2158	rBV4	131338	184000	9.60%	0.906%
82	14.962	2167	2172	2180	rVB3	129448	274805	14.33%	1.353%
83	15.120	2192	2198	2205	rBV5	125081	259010	13.51%	1.275%
84	15.504	2256	2261	2266	rBV6	83029	144980	7.56%	0.714%
85	15.578	2266	2273	2278	rBV5	59787	154739	8.07%	0.762%
86	15.925	2324	2330	2337	rVB2	93082	186929	9.75%	0.920%
87	16.041	2345	2349	2354	rVB	64530	100797	5.26%	0.496%
88	16.224	2375	2379	2386	rVB5	67789	135443	7.07%	0.667%
89	16.364	2396	2402	2415	rVB8	75697	211550	11.04%	1.042%
90	16.474	2415	2420	2429	rVB	98696	199024	10.38%	0.980%

Sum of corrected areas: 20309785

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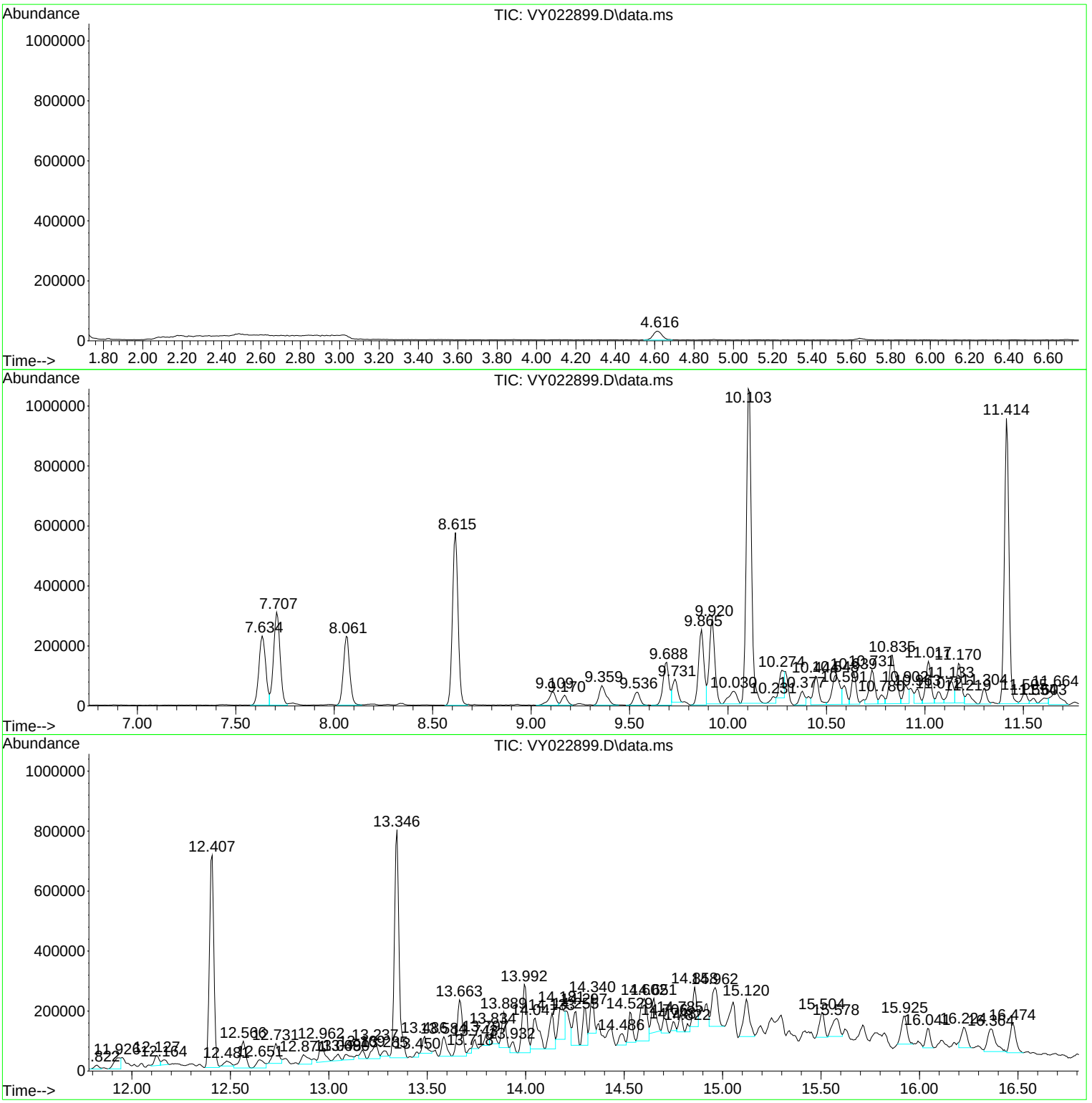
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
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Instrument :
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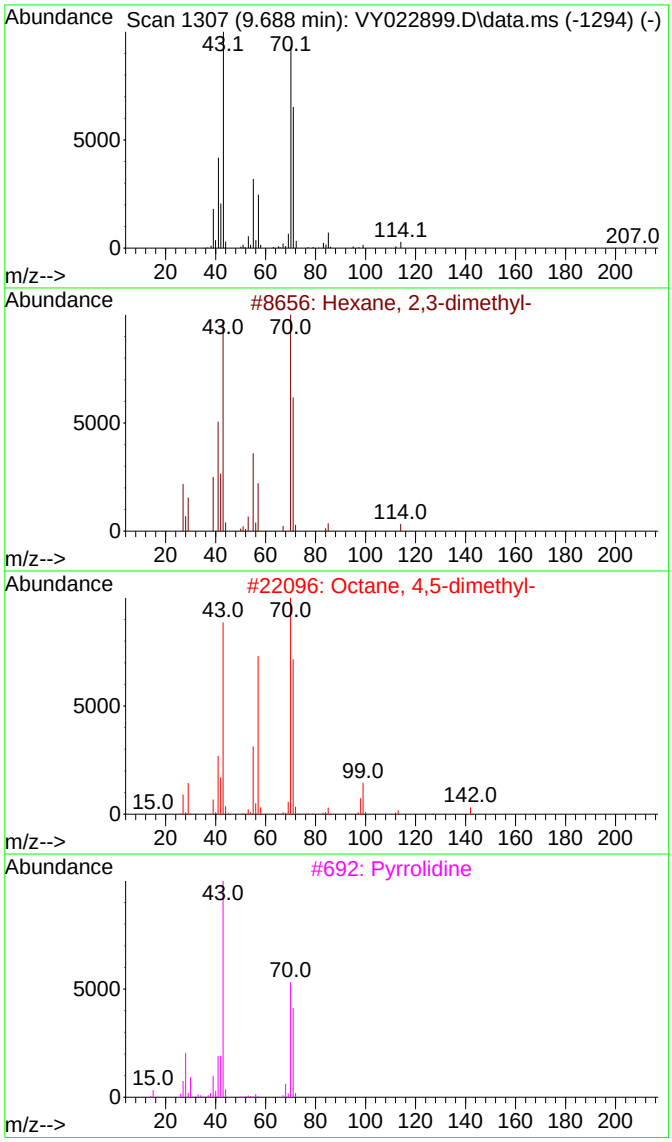
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

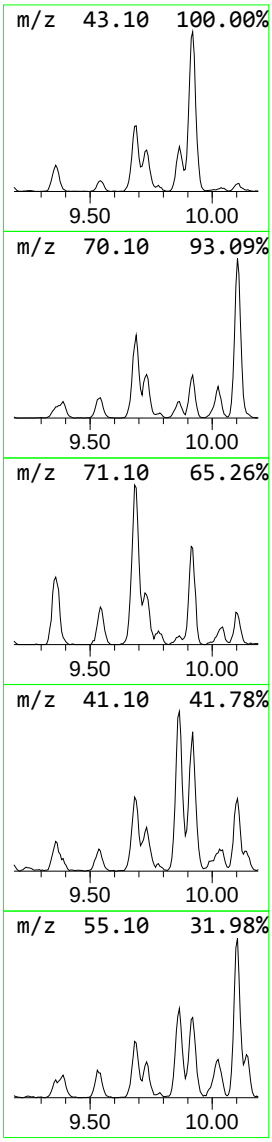
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Hexane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.688	13.70 ug/l	310702	1,4-Difluorobenzene	8.615

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	91
2	Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	83
3	Pyrrolidine	71	C4H9N	000123-75-1	78
4	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	72
5	Heptane, 4-methyl-	114	C8H18	000589-53-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

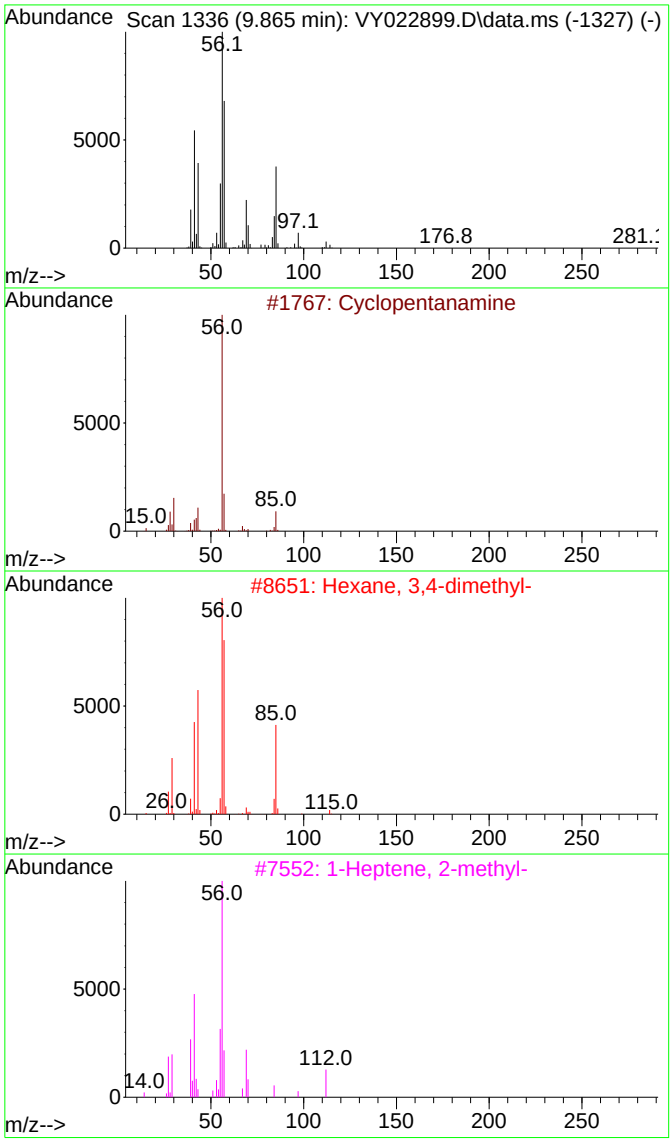
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

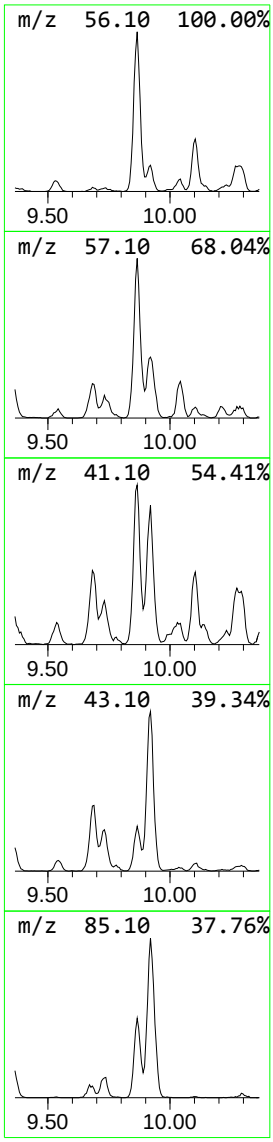
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclopentanamine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.865	21.65 ug/l	490761	1,4-Difluorobenzene	8.615

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanamine	85	C5H11N	001003-03-8	72
2	Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	72
3	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	59
4	2,4-Dimethyl-1-hexene	112	C8H16	016746-87-5	58
5	Hexyl alcohol, chlorodifluoroace...	214	C8H13ClF2O2	1000447-24-1	56





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Instrument :
MSVOA_Y
ClientSampleId :
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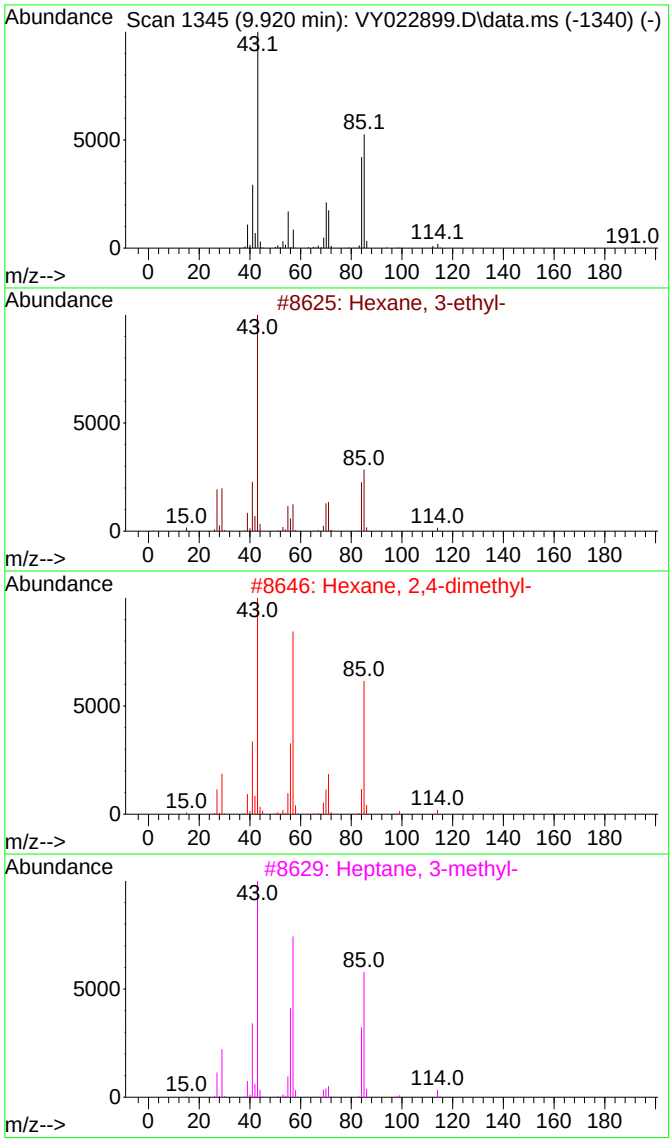
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

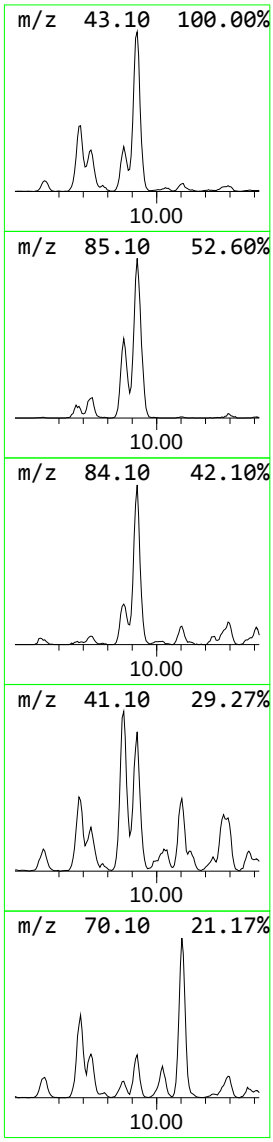
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexane, 3-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.920	24.06 ug/l	545420	1,4-Difluorobenzene	8.615

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-ethyl-	114	C8H18	000619-99-8	91
2	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	52
3	Heptane, 3-methyl-	114	C8H18	000589-81-1	50
4	Nonane, 5-methyl-	142	C10H22	015869-85-9	45
5	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	40





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ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

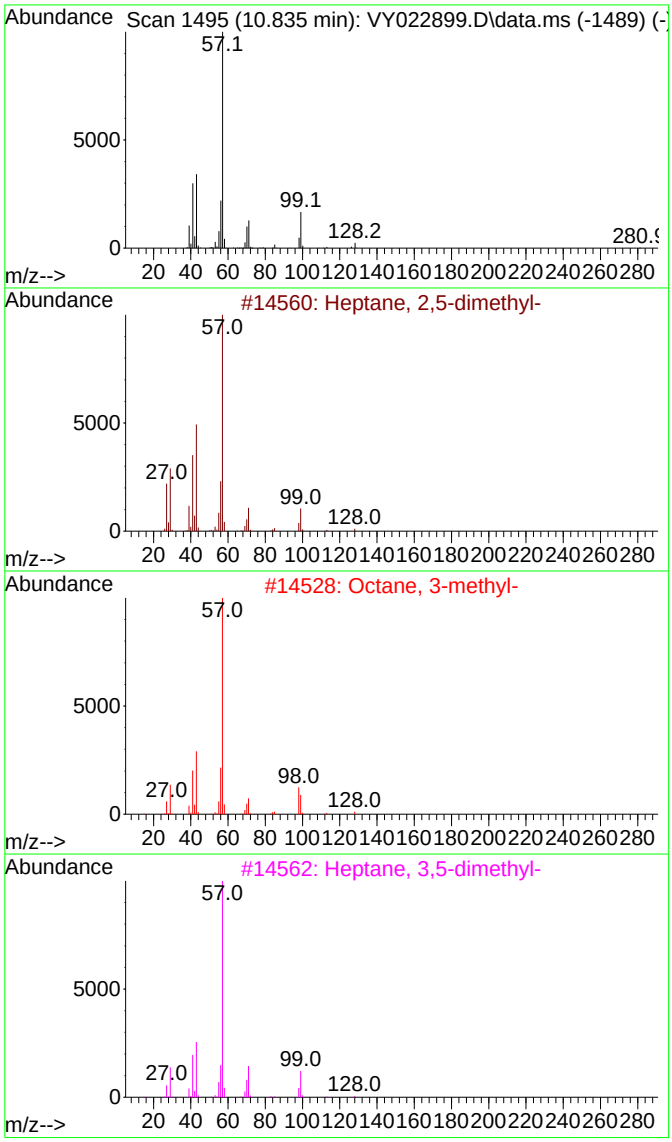
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

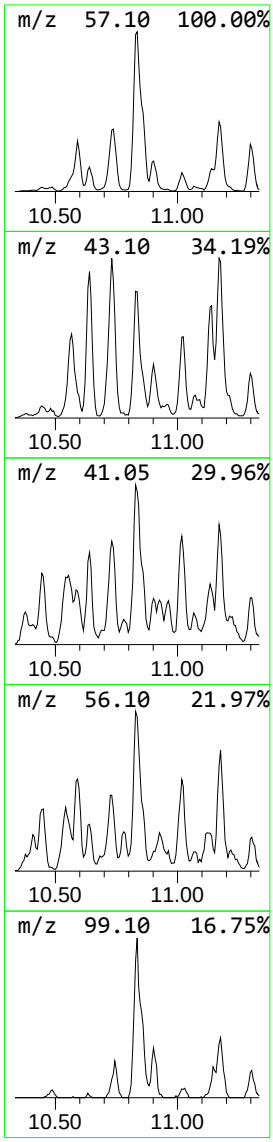
****

Peak Number 4 Heptane, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.835	12.42 ug/l	385300	Chlorobenzene-d5	11.414

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	91
2	Octane, 3-methyl-	128	C9H20	002216-33-3	90
3	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	81
4	Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	47
5	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	45





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Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

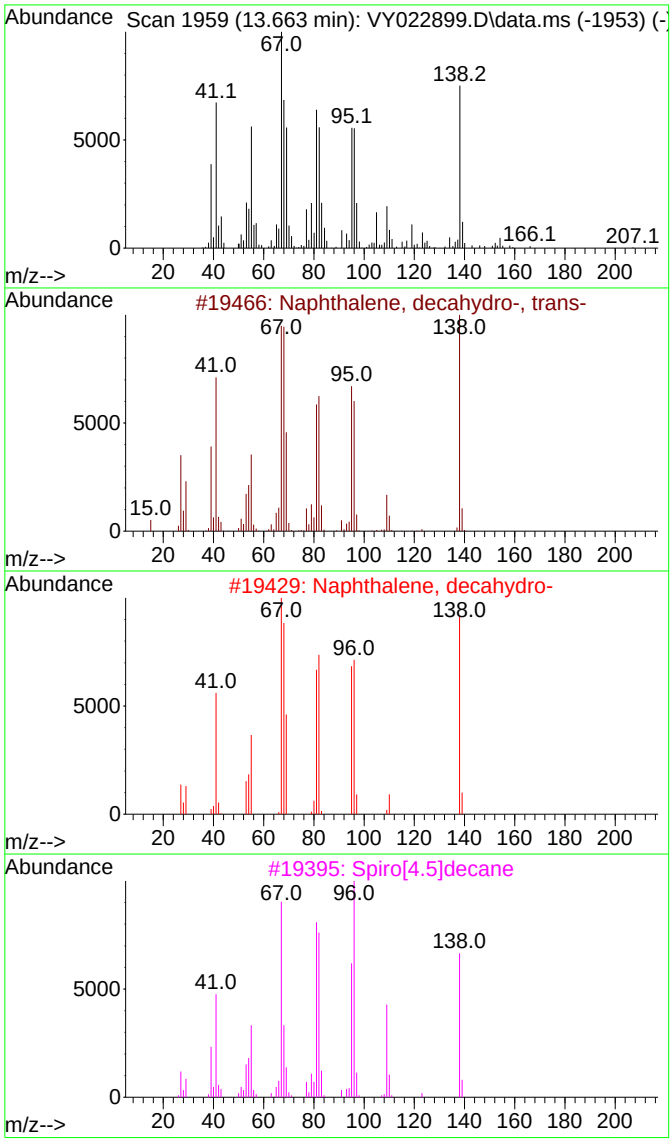
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

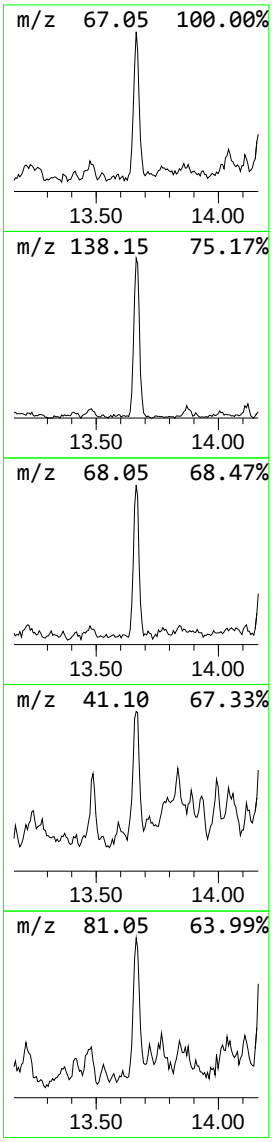
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Naphthalene, decahydro-, tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	13.92 ug/l	348238	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
2	Naphthalene, decahydro-	138	C10H18	000091-17-8	95
3	Spiro[4.5]decane	138	C10H18	000176-63-6	81
4	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	81
5	1,1'-Bicyclopentyl	138	C10H18	001636-39-1	76





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Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

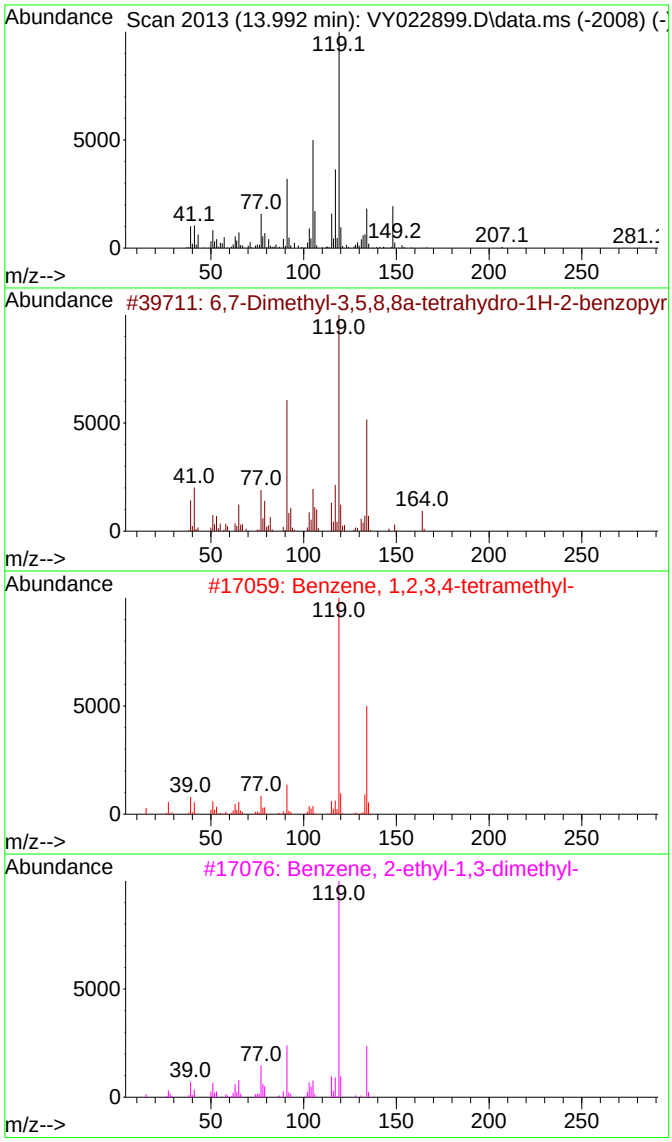
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

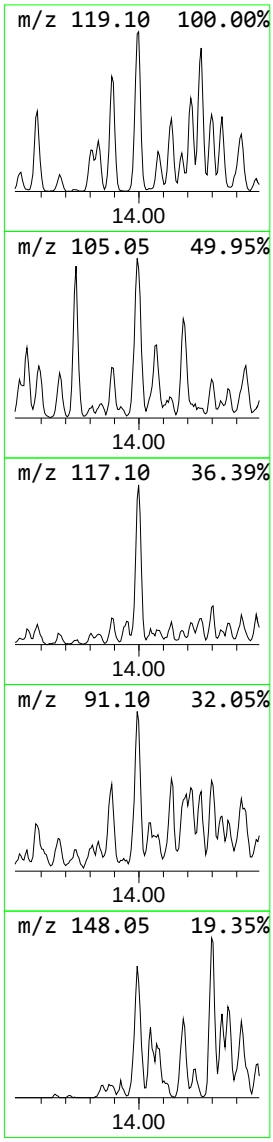
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 6,7-Dimethyl-3,5,8,8a-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.993	15.76 ug/l	394176	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6,7-Dimethyl-3,5,8,8a-tetrahydro...	164	C11H16O	110028-10-9	72
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
4	o-Cymene	134	C10H14	000527-84-4	64
5	p-Cymene	134	C10H14	000099-87-6	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

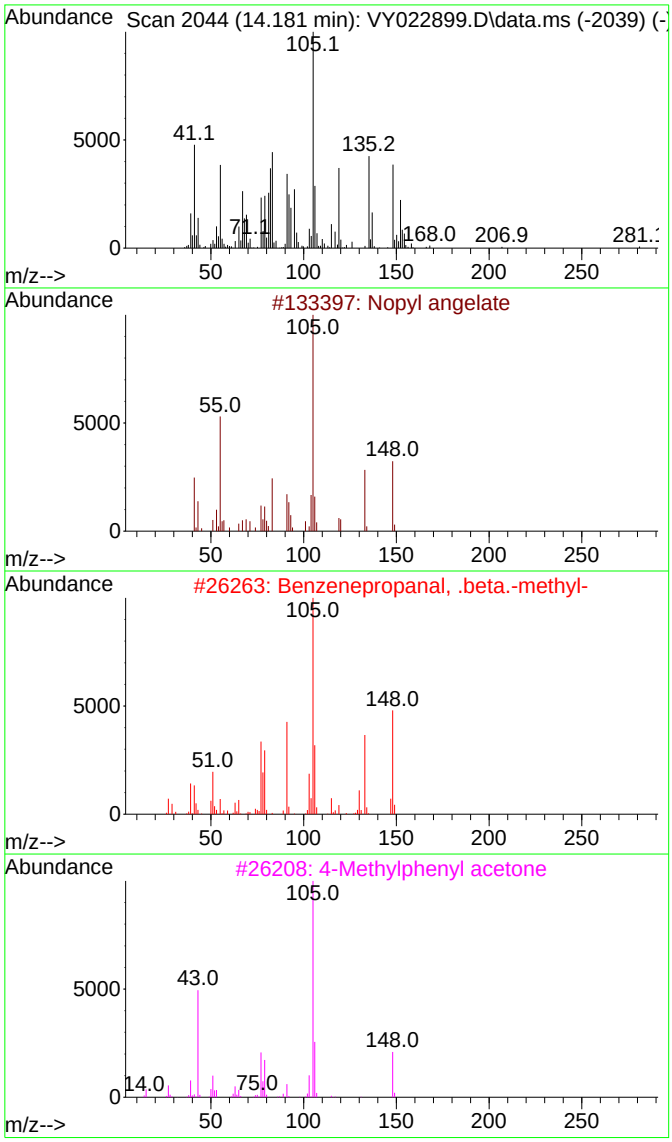
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

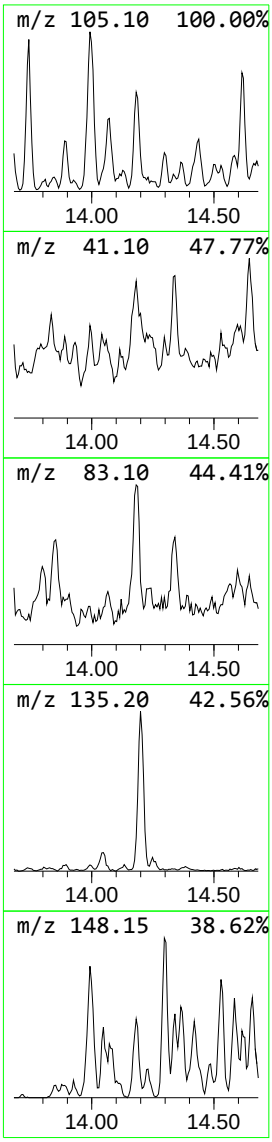
****

Peak Number 7 unknown14.182 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.182	9.51 ug/l	237744	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nopyl angelate	248	C16H24O2	1000383-60-6	32
2	Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	18
3	4-Methylphenyl acetone	148	C10H12O	002096-86-8	14
4	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	14
5	2-Methylindan-2-ol	148	C10H12O	033223-84-6	14





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

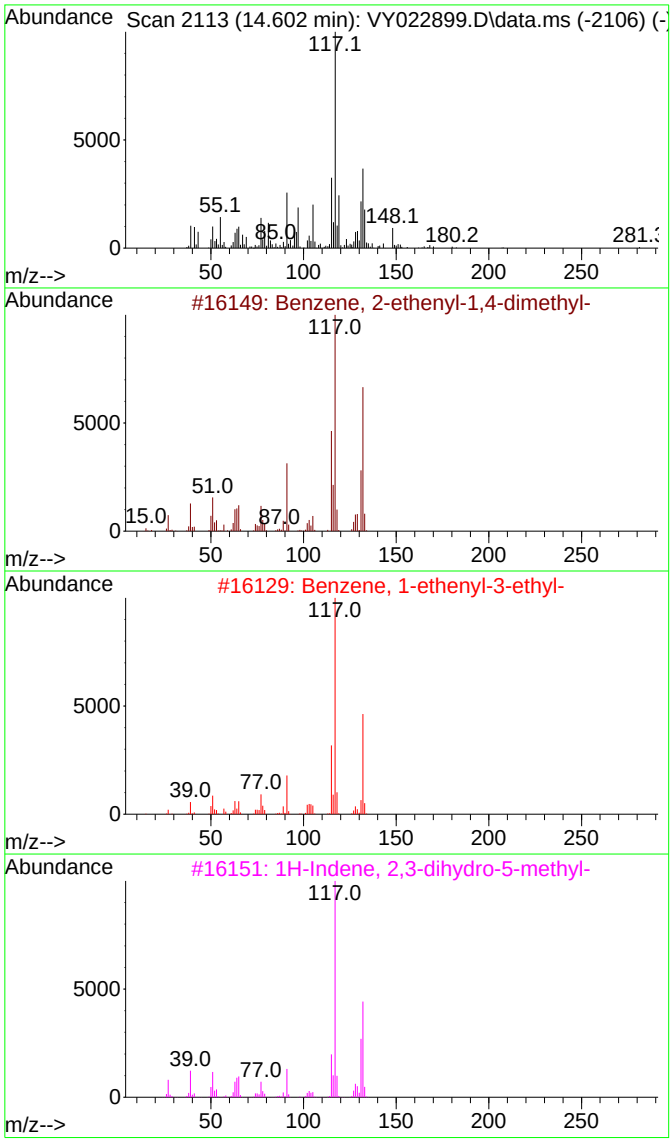
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

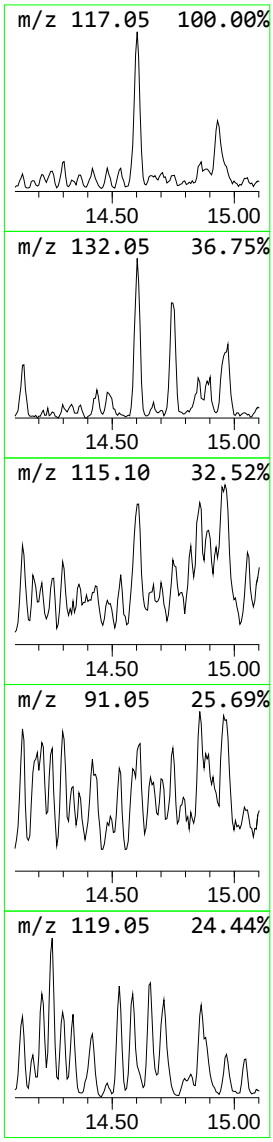
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.602	14.25 ug/l	356504	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	68





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

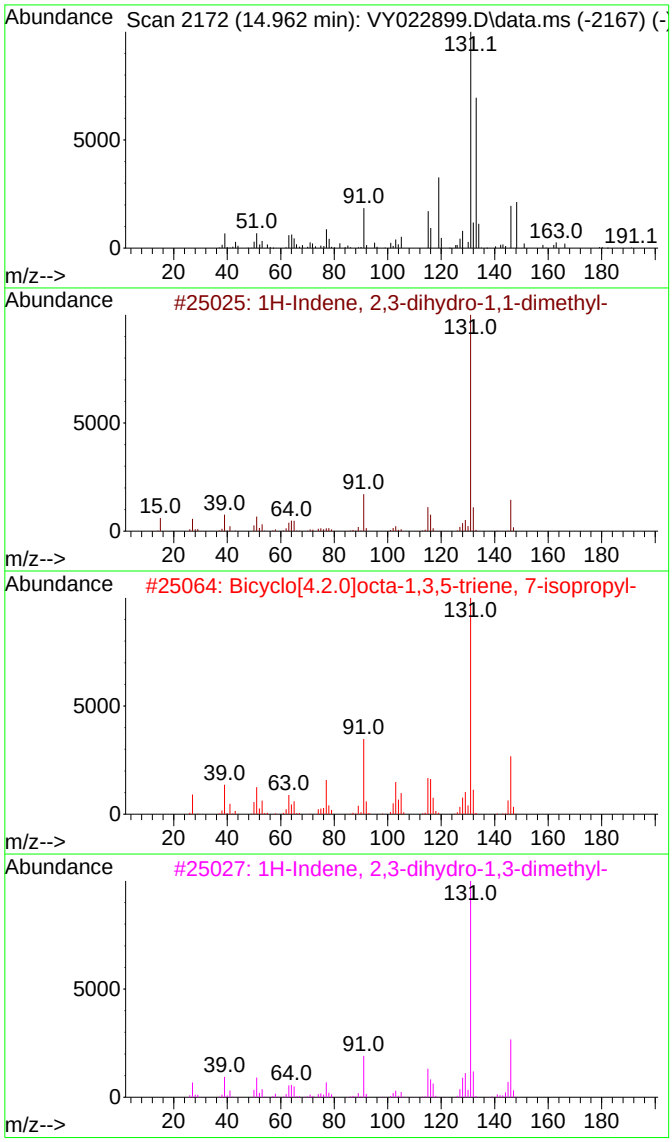
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

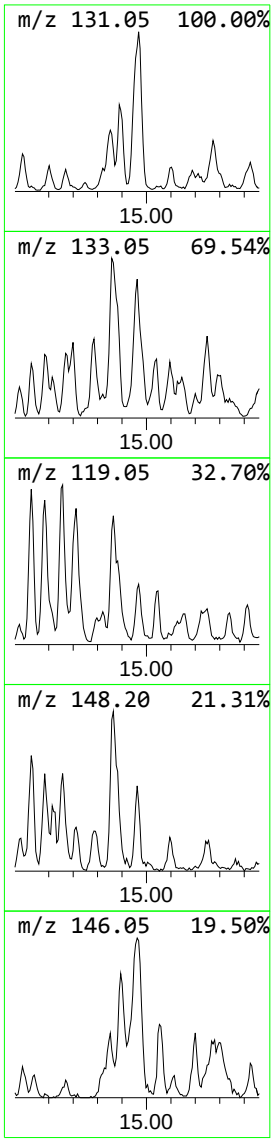
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.962	10.99 ug/l	274805	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	60
2	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	55
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	53
4	Cyclopropane, 1-chloro-1-methyl-...	166	C10H11Cl	1000161-07-9	43
5	2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

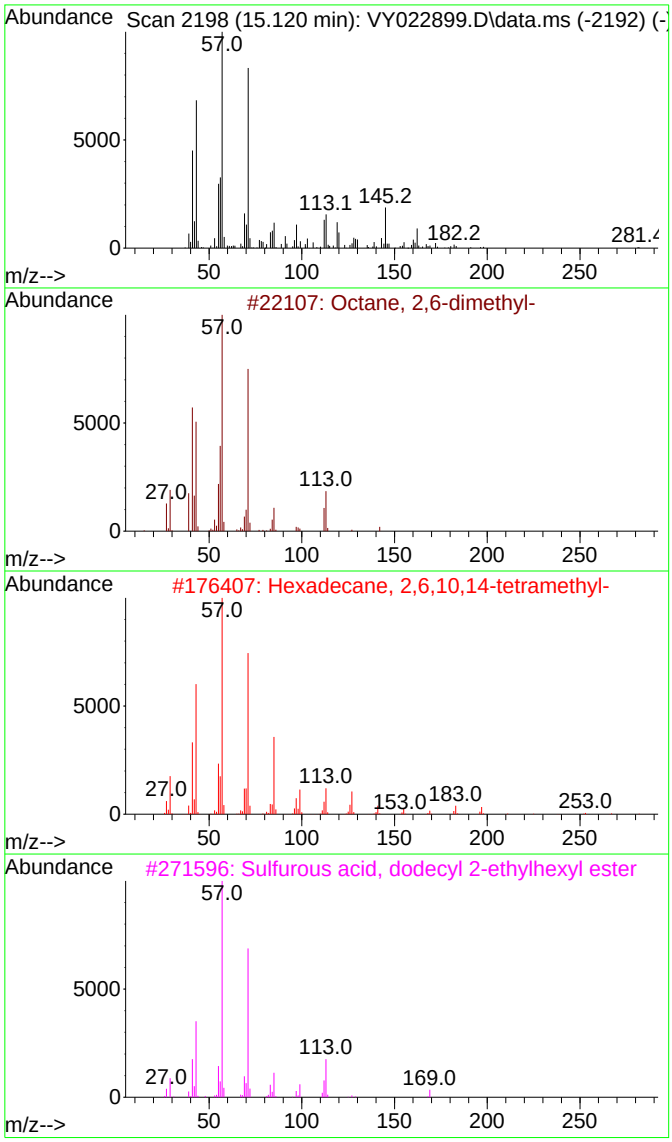
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

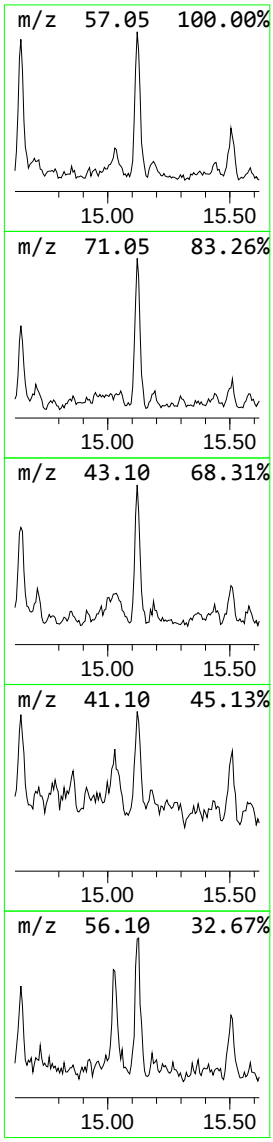
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Octane, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.120	10.36 ug/l	259010	1,4-Dichlorobenzene-d4	13.346

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	58
3	Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	53
4	Octyl tetracosyl ether	467	C32H66O	1000406-39-0	53
5	Tridecane, 7-methyl-	198	C14H30	026730-14-3	52





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070125\
Data File : VY022899.D
Acq On : 01 Jul 2025 15:06
Operator : SY/MD
Sample : Q2458-05
Misc : 6.39g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-66

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062325S.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Hexane, 2,3-dim...	9.688	13.7	ug/l	310702	2	8.615	1133630	50.0
Cyclopentanamine	9.865	21.6	ug/l	490761	2	8.615	1133630	50.0
Hexane, 3-ethyl-	9.920	24.1	ug/l	545420	2	8.615	1133630	50.0
Heptane, 2,5-di...	10.835	12.4	ug/l	385300	3	11.414	1551420	50.0
Naphthalene, de...	13.663	13.9	ug/l	348238	4	13.346	1250490	50.0
6,7-Dimethyl-3,...	13.993	15.8	ug/l	394176	4	13.346	1250490	50.0
unknown14.182	14.182	9.5	ug/l	237744	4	13.346	1250490	50.0
Benzene, 2-ethe...	14.602	14.3	ug/l	356504	4	13.346	1250490	50.0
1H-Indene, 2,3-...	14.962	11.0	ug/l	274805	4	13.346	1250490	50.0
Octane, 2,6-dim...	15.120	10.4	ug/l	259010	4	13.346	1250490	50.0