LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

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_N\methods\82N121422W.M
Title : SW846 8260

Signal : TIC: VN075868.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.253	192	204	222	rBV	1788518	4206936	4.82%	0.681%
2	3.653	262	272	292	rVB2	774431	2209569	2.53%	0.358%
3	4.159	348	358	376	rVB	480147	1354110	1.55%	0.219%
4	4.436	391	405	425	rBV2	330416	1141181	1.31%	0.185%
5	5.289	538	550	553	rBV2	527740	1584455	1.81%	0.257%
6	5.353	554	561	601	rVB	1326161	5481480	6.28%	0.887%
7	5.824	627	641	663	rBV	838999	2906331	3.33%	0.471%
8	6.324	712	726	739	rBV	1062746	3202144	3.67%	0.518%
9	7.088	846	856	868	rVB2	316563	919207	1.05%	0.149%
10	7.230	868	880	889	rBV	599282	1746763	2.00%	0.283%
11	7.336	889	898	920	rVB	1075521	3180071	3.64%	0.515%
12	8.012	1004	1013	1022	rVB	464886	1088499	1.25%	0.176%
13	8.230	1043	1050	1062	rVV2	2107618	6783967	7.77%	1.098%
14	8.335	1062	1068	1078	rVB2	964102	2306356	2.64%	0.373%
15	8.453	1078	1088	1096	rBV	1966599	4338453	4.97%	0.702%
16	8.612	1104	1115	1125	rVB	1534905	3791818	4.34%	0.614%
17	8.771	1125	1142	1152	rBV	7306080	19096627	21.86%	3.092%
18	8.865	1152	1158	1166	rVB	393915	908753	1.04%	0.147%
19	8.971	1166	1176	1190	rVB	1904627	3982778	4.56%	0.645%
20	9.106	1190	1199	1205	rBV2	977757	2172382	2.49%	0.352%
21	9.600	1267	1283	1289	rBV2	3004383	7799224	8.93%	1.263%
22	9.647	1289	1291	1299	rVV	1047394	1594855	1.83%	0.258%
23	9.729	1299	1305	1310	rVV	476028	1046831	1.20%	0.169%
24	10.018	1344	1354	1363	rBV	3362447	6721872	7.70%	1.088%
25	10.153	1363	1377	1383	rBV	5030590	10705681	12.26%	1.733%
26	10.206	1383	1386	1391	rVV	650149	949839	1.09%	0.154%
27	10.347	1400	1410	1422	rBV	1038430	2469644	2.83%	0.400%
28	10.500	1430	1436	1442	rVB2	911351	1679386	1.92%	0.272%
29	10.571	1442	1448	1452	rBV	1778402	3280911	3.76%	0.531%
30	10.635	1452	1459	1467	rVB2	48052415	87339035	100.00%	14.140%
31	10.747	1472	1478	1485	rVB2	1950907	3672132	4.20%	0.595%
32	11.012	1514	1523	1531	rBV4	727534	1669518	1.91%	0.270%
33	11.424	1588	1593	1597	rVV	870687	1618359	1.85%	0.262%
34	11.476	1597	1602	1607	rVB	576077	1066370	1.22%	0.173%
35	11.647	1623	1631	1634	rBV2	616794	1216820	1.39%	0.197%
36	11.747	1644	1648	1654	rVB	595055	916221	1.05%	0.148%

LSC Area Percent Report

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ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

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_N\methods\82N121422W.M
Title : SW846 8260

37	11.871	1662	1669	1674	rBV	789635	1371305	1.57%	0.222%
38	11.971	1678	1686	1695	rVV	13125088	22256202	25.48%	3.603%
39	12.071	1695	1703	1715	rVV2	43307941	83042114	95.08%	13.445%
40	12.159	1715	1718	1724	rVB2	570636	891248	1.02%	0.144%
41	12.400	1747	1759	1768	rBV	18941985	33025696	37.81%	5.347%
42	12.588	1785	1791	1793	rBV2	485215	912872	1.05%	0.148%
43	12.623	1794	1797	1801	rVV2	617936	987903	1.13%	0.160%
44	12.700	1805	1810	1815	rVV	1420699	2187769	2.50%	0.354%
45	12.853	1831	1836	1840	rBV	661694	1119562	1.28%	0.181%
46	13.041	1860	1868	1872	rBV	4586487	7594520	8.70%	1.230%
47	13.100	1872	1878	1882	rVV	15344527	25880439	29.63%	4.190%
48	13.135	1882	1884	1887	rVV	6859623	9098591	10.42%	1.473%
49	13.176	1887	1891	1897	rVB	7492024	12582226	14.41%	2.037%
50	13.341	1906	1919	1925	rBV	6361903	11205460	12.83%	1.814%
51	13.417	1925	1932	1937	rVB3	424857	1162310	1.33%	0.188%
52	13.488	1937	1944	1954	rBV	26087795	40353782	46.20%	6.533%
53	13.617	1957	1966	1971	rBV2	1065615	2001301	2.29%	0.324%
54	13.676	1971	1976	1981	rBV2	1887344	3058839	3.50%	0.495%
55	13.729	1981	1985	1991	rVB2	738665	1303534	1.49%	0.211%
56	13.823	1991	2001	2014	rBV	6071193	11448995	13.11%	1.854%
57	13.953	2014	2023	2025	rBV3	1485562	2700774	3.09%	0.437%
58	13.988	2025	2029	2033	rVV2	4696943	8530907	9.77%	1.381%
59	14.023	2033	2035	2046	rVB4	4138507	8956726	10.26%	1.450%
60	14.117	2046	2051	2057	rBV3	1836275	3581796	4.10%	0.580%
61	14.188	2057	2063	2068	rBV	1917130	2889149	3.31%	0.468%
62	14.247	2068	2073	2075	rBV	2395016	3487333	3.99%	0.565%
63	14.276	2075	2078	2083	rVB	2404828	3563696	4.08%	0.577%
64	14.335	2083	2088	2094	rVB	3881921	5866730	6.72%	0.950%
65	14.447	2101	2107	2112	rVB2	3125300	5333438	6.11%	0.863%
66	14.576	2124	2129	2134	rBV2	1236739	1884376	2.16%	0.305%
67	14.629	2134	2138	2140	rBV	1132141	1678924	1.92%	0.272%
68	14.659	2140	2143	2146	rVB	1160470	1363429	1.56%	0.221%
69	14.694	2146	2149	2153	rVB	2409522	3205169	3.67%	0.519%
70	14.735	2153	2156	2160	rVB	1384776	1759534	2.01%	0.285%
71	14.776	2160	2163	2171	rVB3	909816	1927586	2.21%	0.312%
72	14.882	2171	2181	2185	rBV2	1666226	3611202	4.13%	0.585%
73	14.935	2185	2190	2193	rBV3	2619430	5362944	6.14%	0.868%
74	15.059	2201	2211	2216	rBV4	4439934	12013453	13.75%	1.945%
75	15.106	2216	2219	2229	rVB3	1061638	2122586	2.43%	0.344%
76	15.217	2232	2238	2244	rVB	2677736	4809619	5.51%	0.779%

LSC Area Percent Report

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Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

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_N\methods\82N121422W.M
Title : SW846 8260

77	15.282	2244	2249	2251	rBV5	668856	1085963	1.24%	0.176%
78	15.323	2251	2256	2261	rBV3	1585194	2911807	3.33%	0.471%
79	15.447	2269	2277	2283	rVB3	2025701	4860274	5.56%	0.787%
80	15.606	2297	2304	2305	rBV3	718606	1232515	1.41%	0.200%
81	15.641	2306	2310	2315	rVB	1368355	2381383	2.73%	0.386%
82	15.700	2315	2320	2325	rBV	1732891	3424131	3.92%	0.554%
83	15.753	2325	2329	2332	rBV2	982644	1512763	1.73%	0.245%
84	15.835	2340	2343	2354	rVB4	1478131	3294123	3.77%	0.533%
85	15.953	2354	2363	2370	rBV3	1658215	4102313	4.70%	0.664%
86	16.129	2381	2393	2394	rVV3	1151990	2997446	3.43%	0.485%
87	16.170	2395	2400	2410	rVV	4056051	8824478	10.10%	1.429%
88	16.258	2411	2415	2421	rVV2	638498	1517573	1.74%	0.246%
89	16.347	2422	2430	2444	rVB3	1151144	4346395	4.98%	0.704%
90	16.506	2447	2457	2473	rBV2	2148139	6259147	7.17%	1.013%
91	16.717	2482	2493	2499	rBV2	2659583	6625911	7.59%	1.073%

Sum of corrected areas: 617658839

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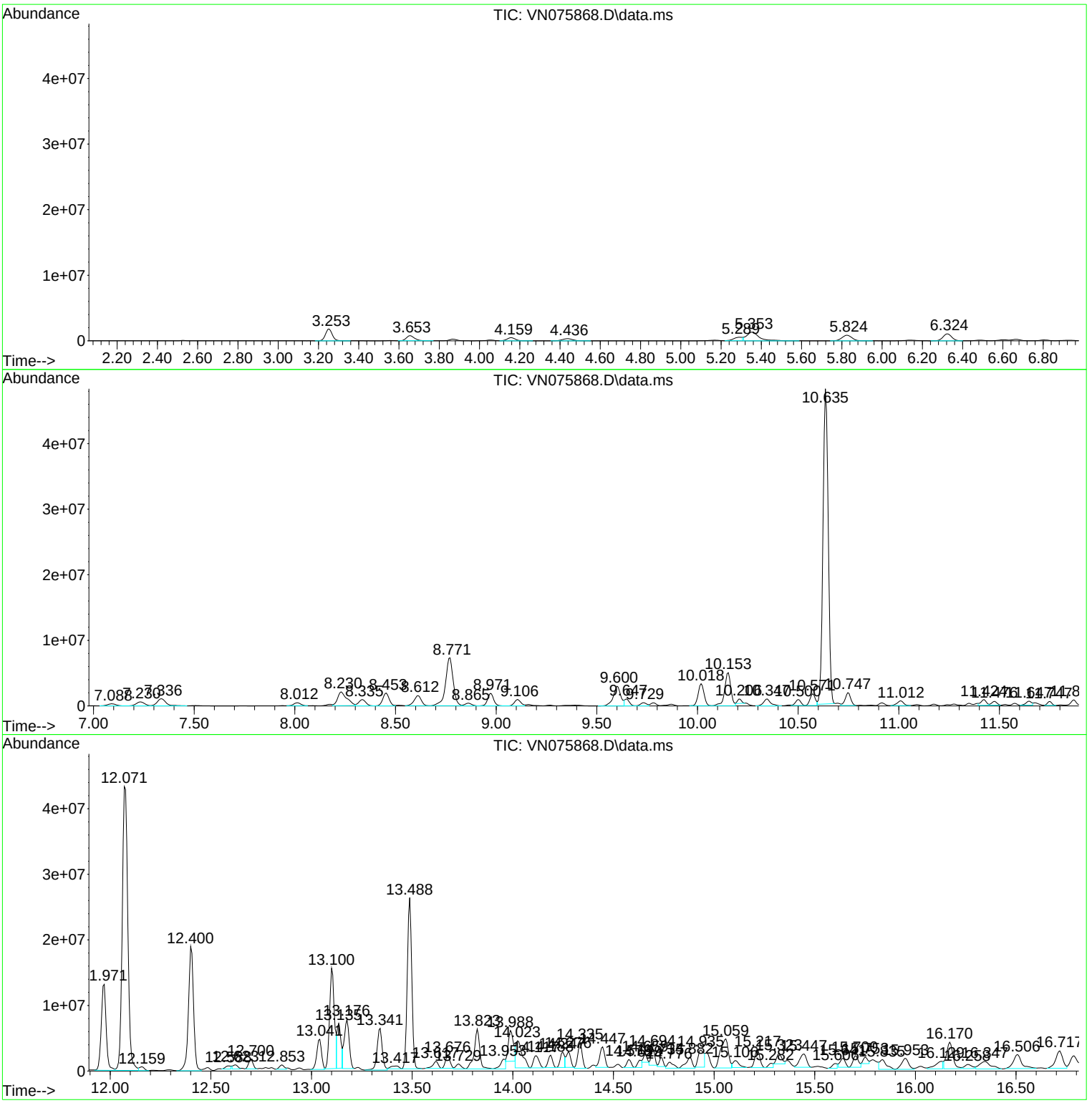
Instrument :
MSVOA_N
ClientSampleId :
5922

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
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Instrument :
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5922

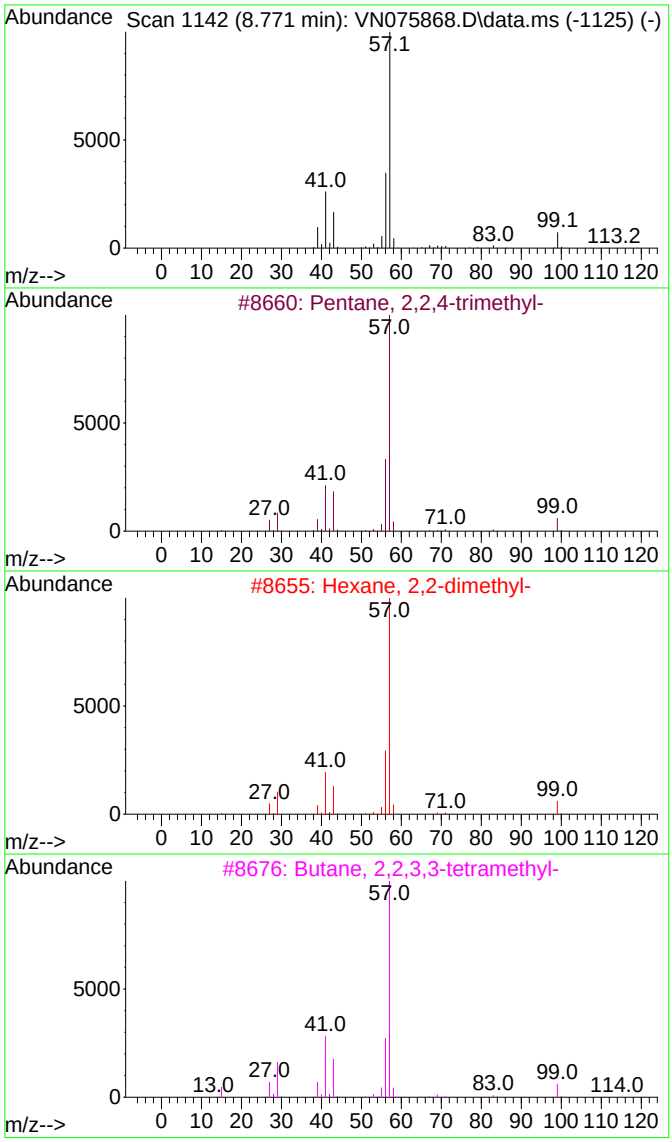
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

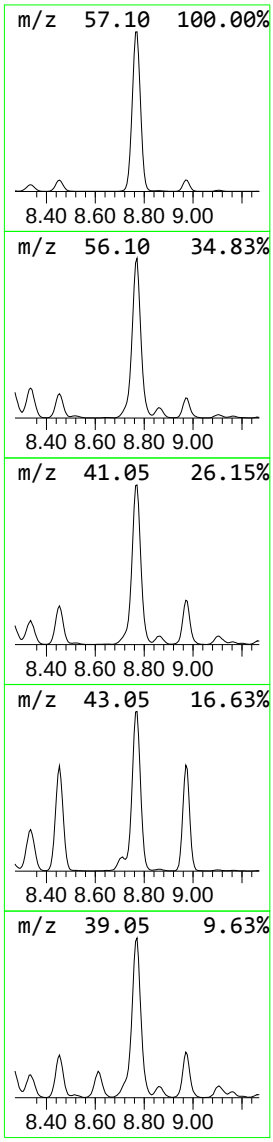
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.771	439.53 ug/l	19096600	1,4-Difluorobenzene	9.106

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
3	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	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	59





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Operator : JC\MD
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Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

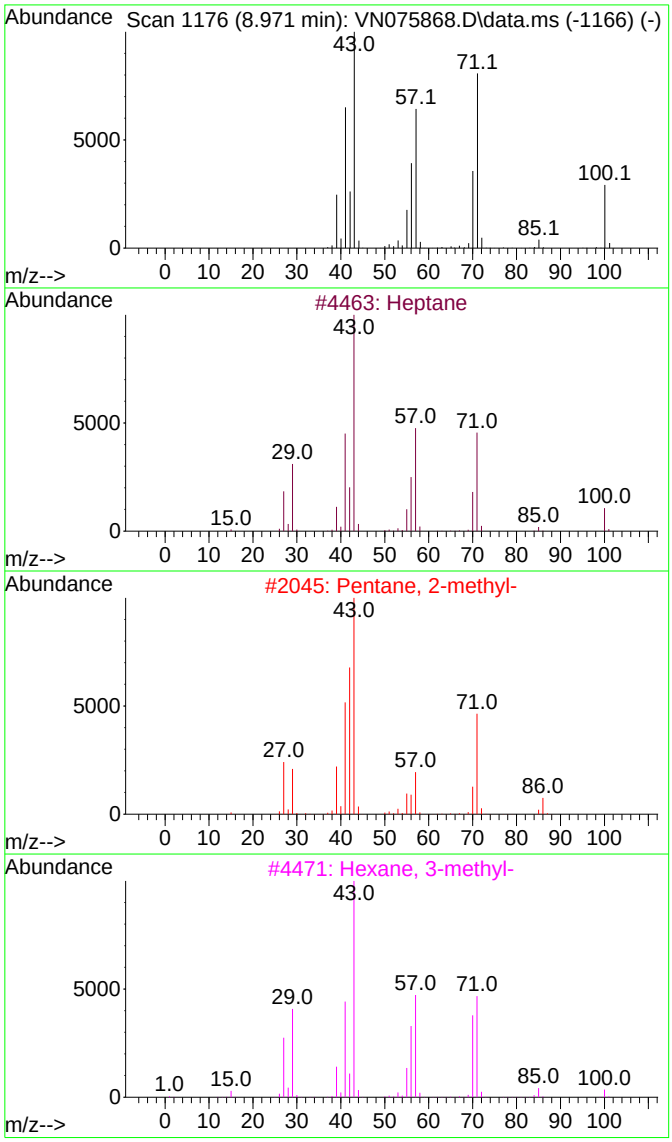
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

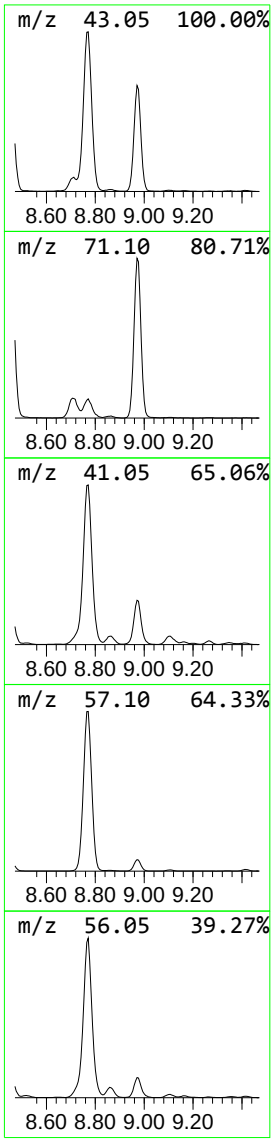
****

Peak Number 2 Heptane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.971	91.67 ug/l	3982780	1,4-Difluorobenzene	9.106

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	50
4	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	47
5	3-Hexanone	100	C6H12O	000589-38-8	43





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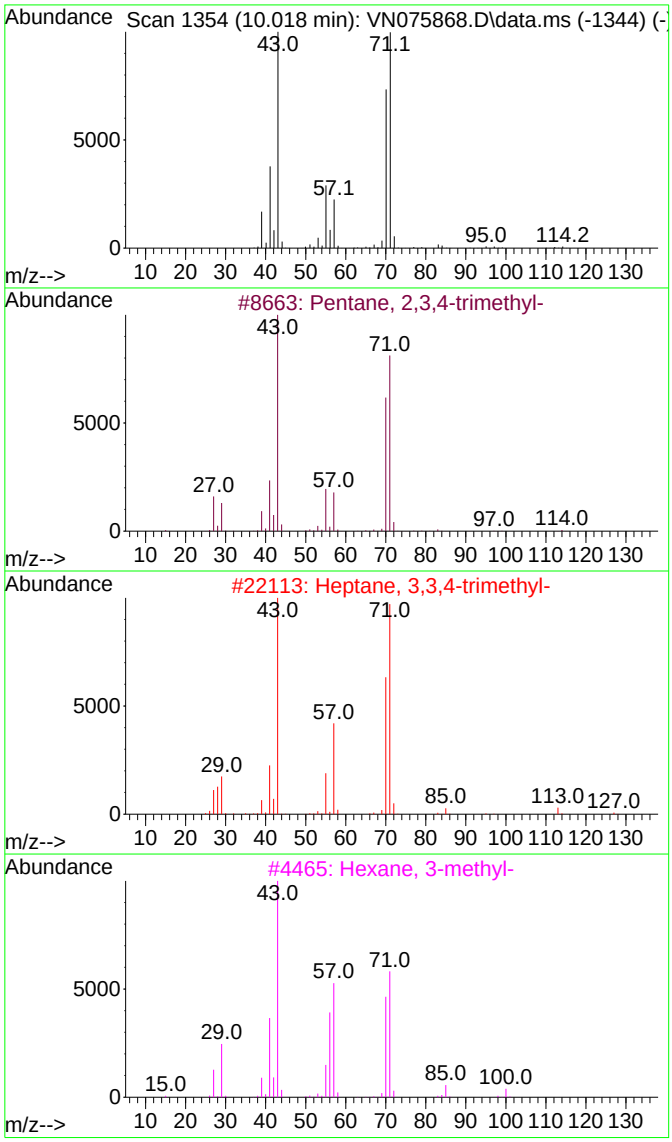
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

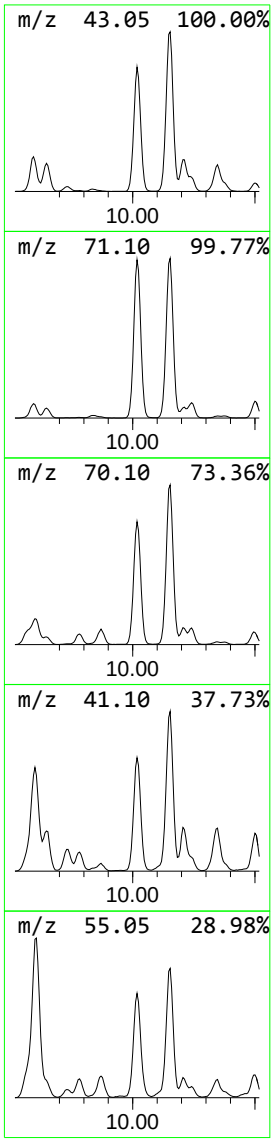
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.018	154.71 ug/l	6721870	1,4-Difluorobenzene	9.106

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	90
2	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	83
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	78
4	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
5	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74





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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

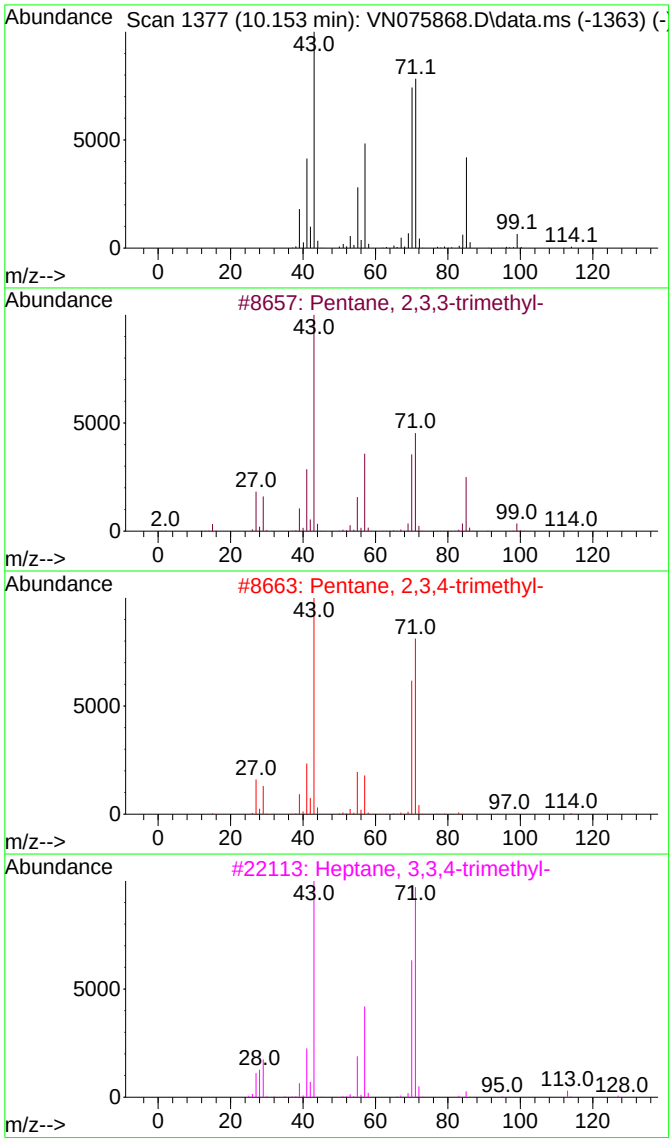
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

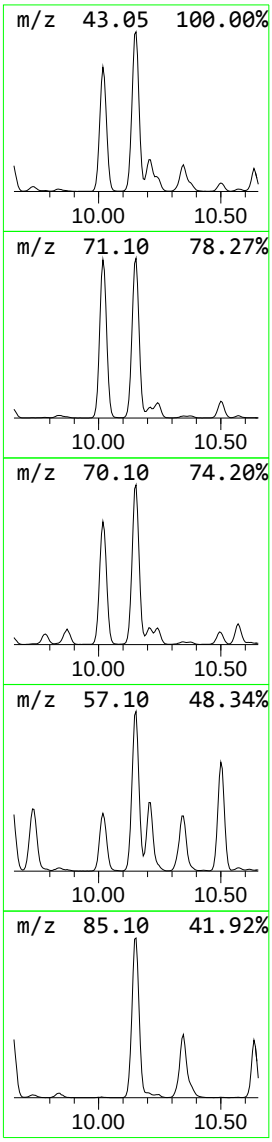
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.153	246.40 ug/l	10705700	1,4-Difluorobenzene	9.106

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	72
3	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	64





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Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

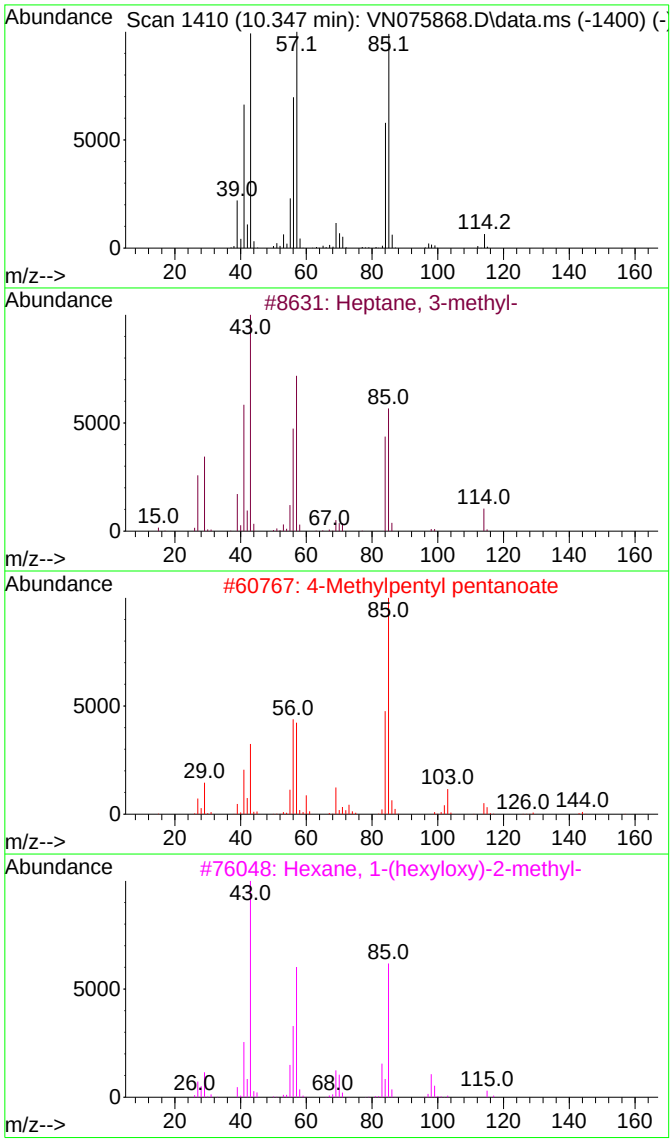
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

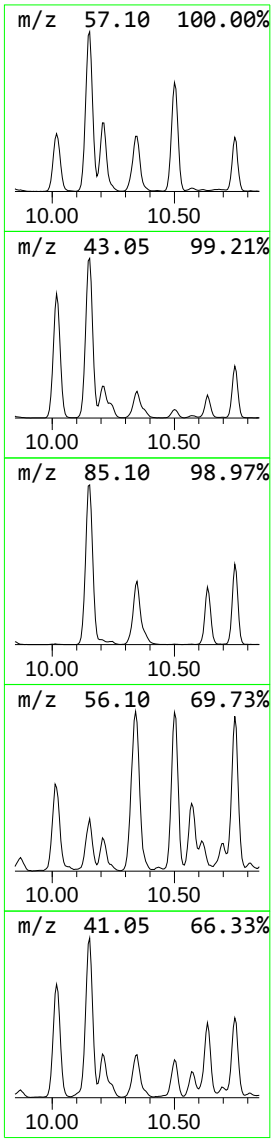
****

Peak Number 5 Heptane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.347	56.84 ug/l	2469640	1,4-Difluorobenzene	9.106

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2	4-Methylpentyl pentanoate	186	C11H22O2	035852-47-2	72
3	Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	64
4	Nonane, 5-methyl-	142	C10H22	015869-85-9	64
5	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	64





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Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

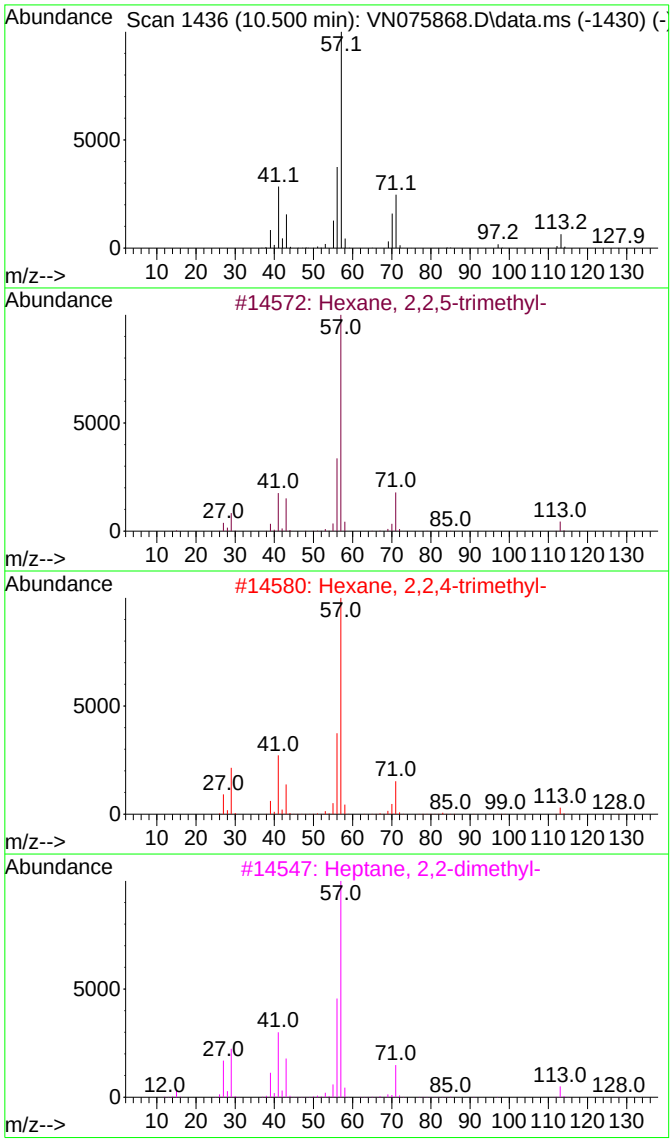
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

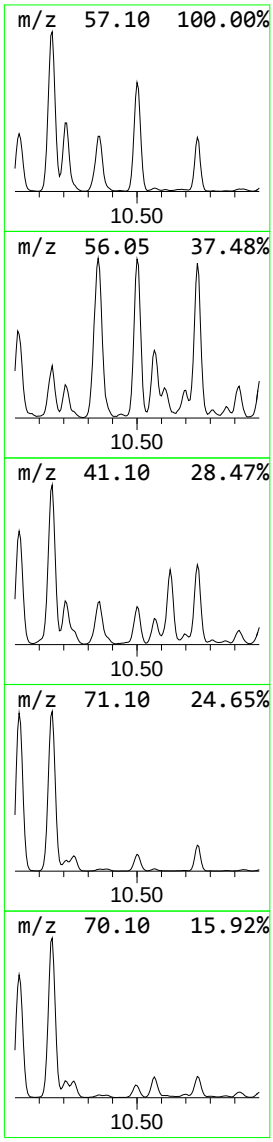
****

Peak Number 6 Hexane, 2,2,5-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.500	61.23 ug/l	1679390	Chlorobenzene-d5	11.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	72
2	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
3	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	59
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
5	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	50





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Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

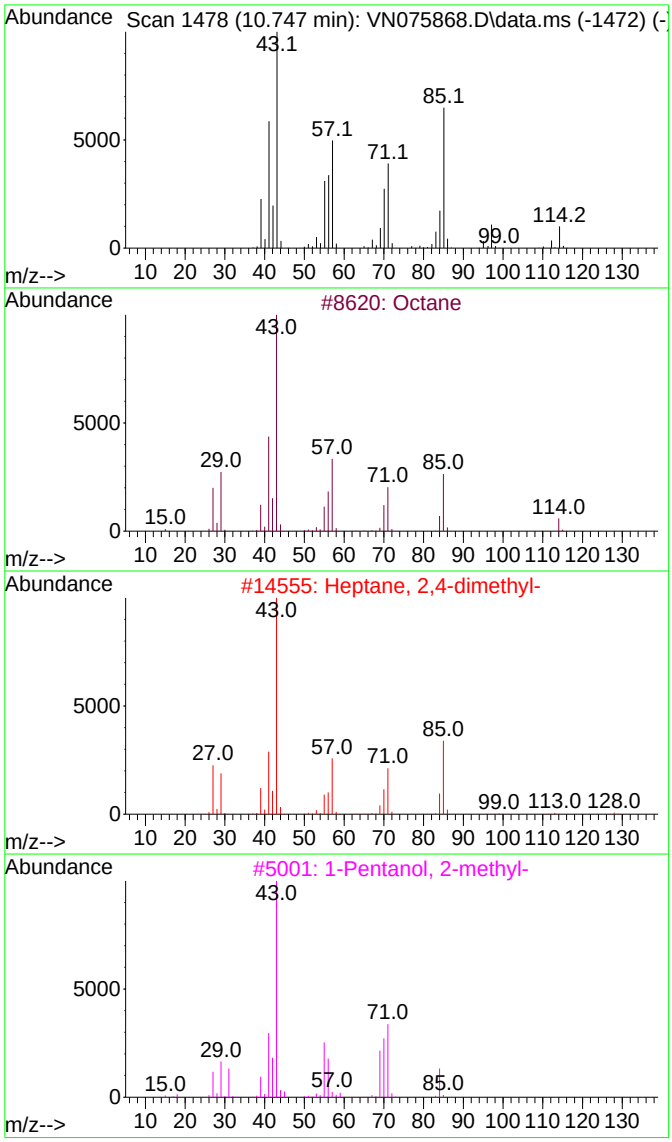
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

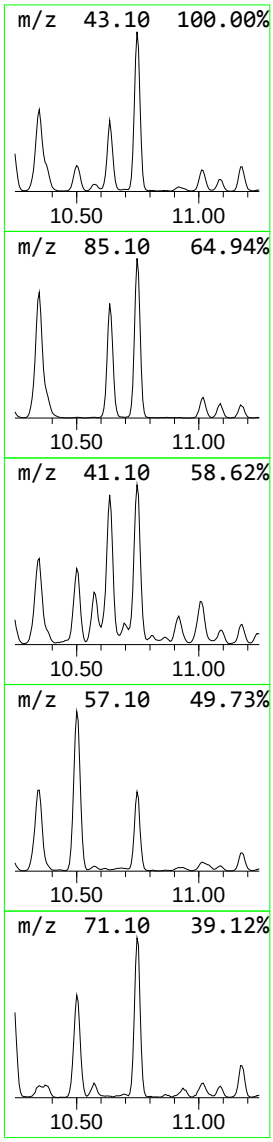
****

Peak Number 7 Octane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.747	133.89 ug/l	3672130	Chlorobenzene-d5	11.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	87
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3	1-Pentanol, 2-methyl-	102	C6H14O	000105-30-6	43
4	Muscimol	114	C4H6N2O2	002763-96-4	43
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	43





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Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

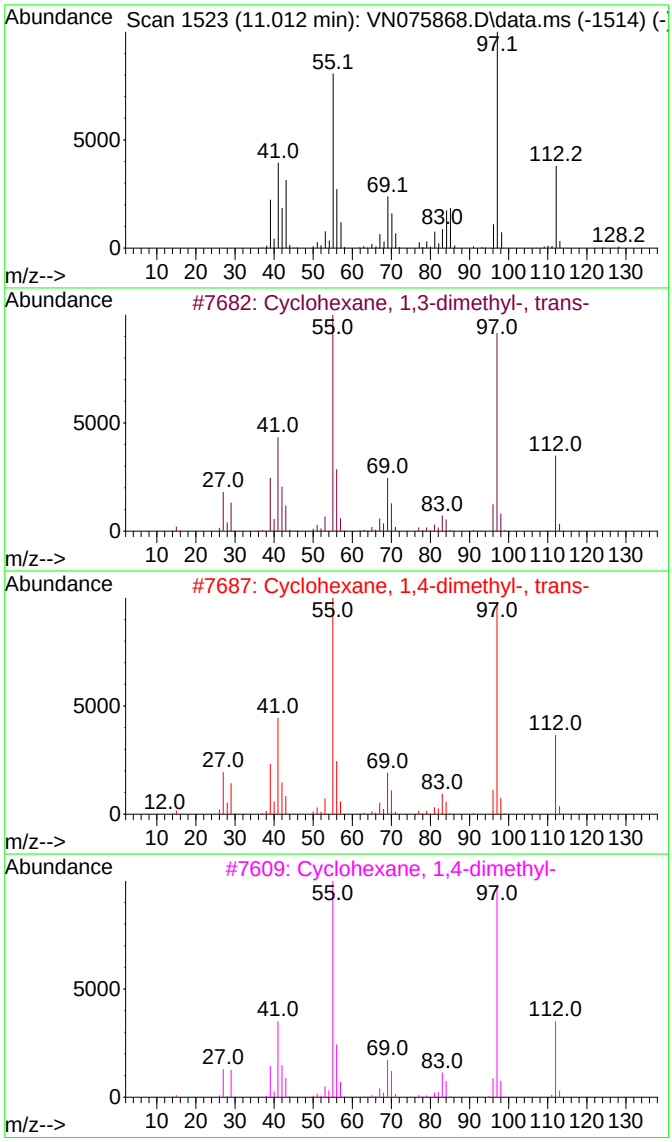
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

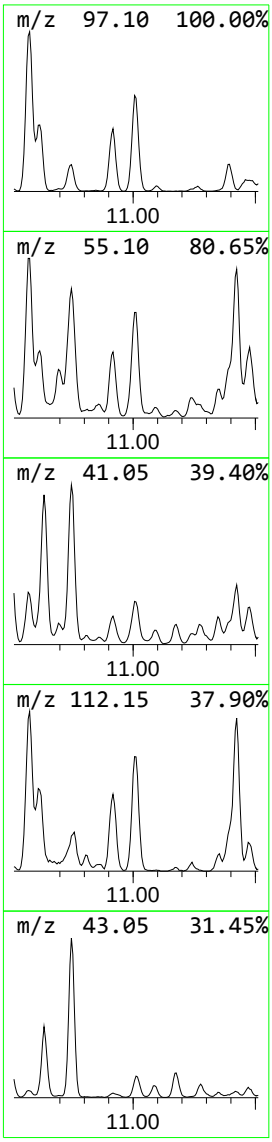
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.012	60.87 ug/l	1669520	Chlorobenzene-d5	11.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	93
2	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	90
3	Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	87
4	Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	81
5	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	80





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Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

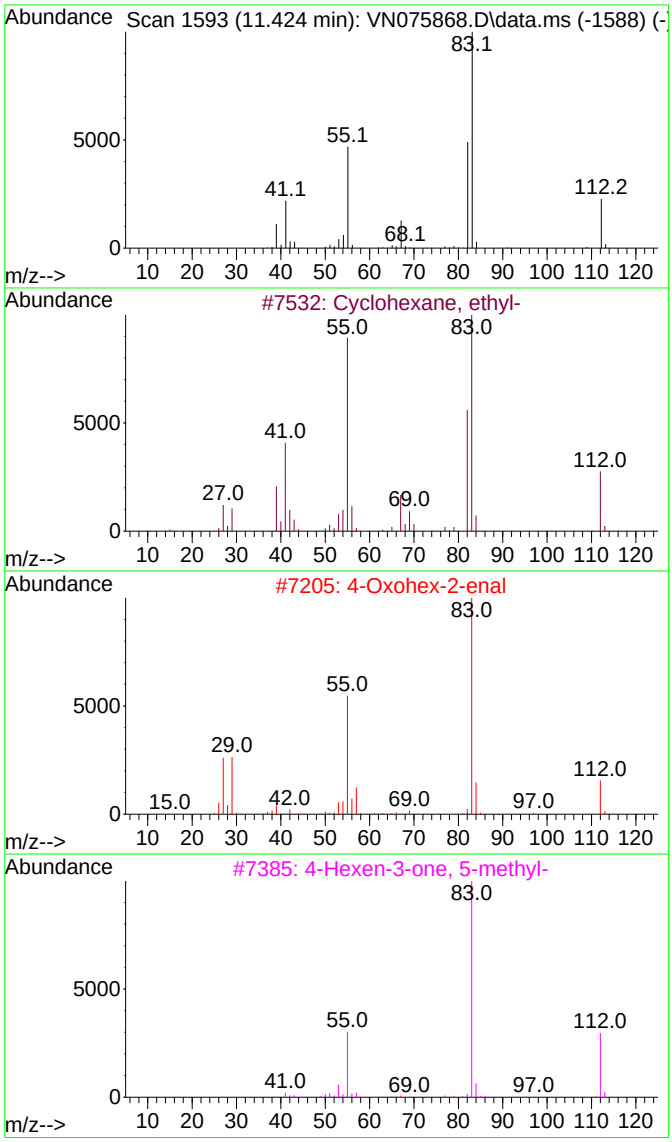
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

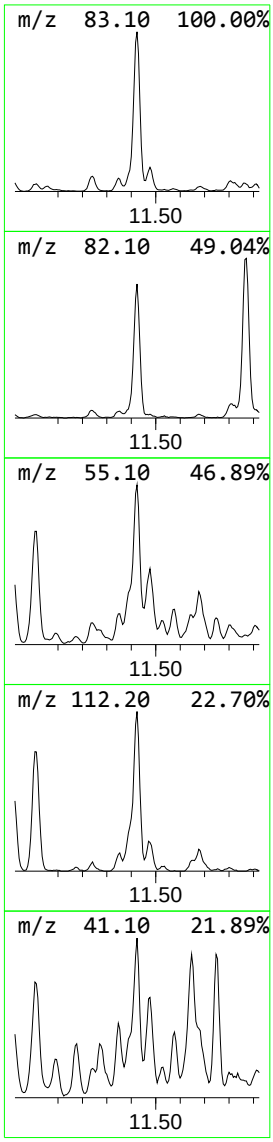
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Cyclohexane, ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.424	59.01 ug/l	1618360	Chlorobenzene-d5	11.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	83
2	4-Oxohex-2-enal	112	C6H8O2	020697-55-6	58
3	4-Hexen-3-one, 5-methyl-	112	C7H12O	013905-10-7	47
4	Cyclohexane, decyl-	224	C16H32	001795-16-0	39
5	Cyclohexane, bromo-	162	C6H11Br	000108-85-0	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075868.D
Acq On : 19 Dec 2022 13:58
Operator : JC\MD
Sample : N6066-21 10X
Misc : 1.10g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5922

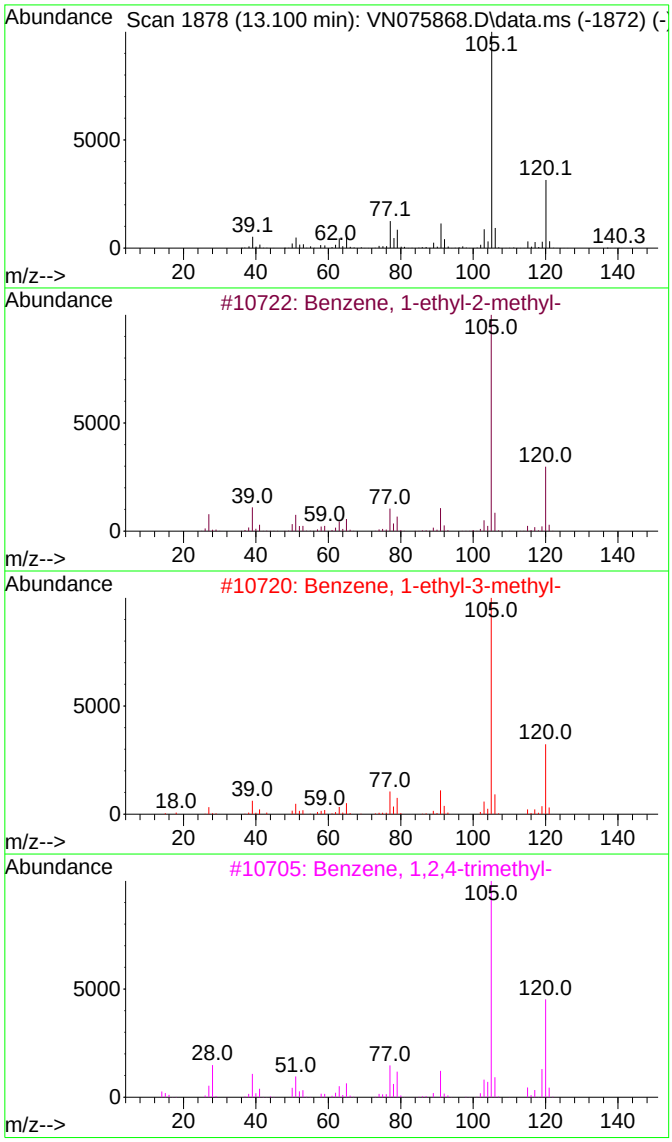
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

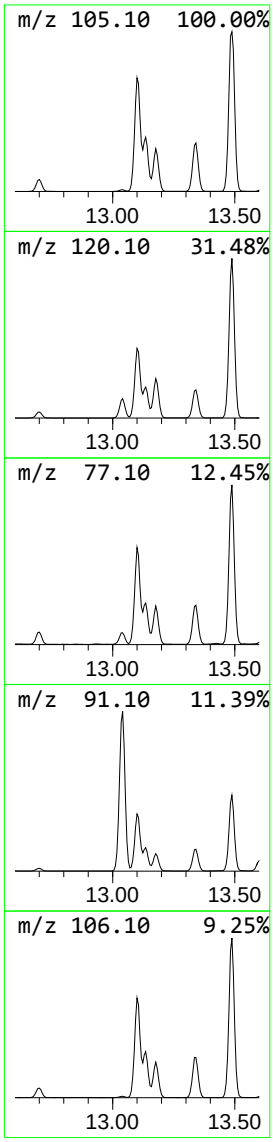
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	113.03 ug/l	25880400	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Tentatively Identified Compound (LSC) summary

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Data File : VN075868.D
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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
Pentane, 2,2,4-...	8.771	439.5	ug/l	19096600	2	9.106	2172380	50.0
Heptane	8.971	91.7	ug/l	3982780	2	9.106	2172380	50.0
Pentane, 2,3,4-...	10.018	154.7	ug/l	6721870	2	9.106	2172380	50.0
Pentane, 2,3,3-...	10.153	246.4	ug/l	10705700	2	9.106	2172380	50.0
Heptane, 3-methyl-	10.347	56.8	ug/l	2469640	2	9.106	2172380	50.0
Hexane, 2,2,5-t...	10.500	61.2	ug/l	1679390	3	11.871	1371310	50.0
Octane	10.747	133.9	ug/l	3672130	3	11.871	1371310	50.0
Cyclohexane, 1,...	11.012	60.9	ug/l	1669520	3	11.871	1371310	50.0
Cyclohexane, et...	11.424	59.0	ug/l	1618360	3	11.871	1371310	50.0
Benzene, 1-ethy...	13.100	113.0	ug/l	25880400	4	13.794	11449000	50.0