LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

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\82N091024W.M
Title : SW846 8260

Signal : TIC: VN083967.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.318	565	572	586	rVB	126030	419620	1.03%	0.137%
2	7.212	882	894	904	rBV2	211005	602495	1.49%	0.197%
3	8.218	1058	1065	1075	rVV2	735577	1838361	4.53%	0.601%
4	8.324	1075	1083	1093	rVB	363319	921094	2.27%	0.301%
5	8.441	1093	1103	1112	rBV	601238	1357155	3.35%	0.443%
6	8.753	1139	1156	1168	rBV	3518367	8874179	21.88%	2.899%
7	8.965	1181	1192	1205	rVB	424775	975544	2.40%	0.319%
8	9.100	1205	1215	1221	rBV2	550975	1248852	3.08%	0.408%
9	9.588	1285	1298	1303	rBV2	1398407	3207202	7.91%	1.048%
10	9.641	1303	1307	1314	rVB	971921	1769383	4.36%	0.578%
11	9.724	1314	1321	1326	rBV	301180	624060	1.54%	0.204%
12	10.012	1362	1370	1380	rBV	3072473	6025550	14.85%	1.969%
13	10.141	1380	1392	1398	rBV	4130003	8400236	20.71%	2.744%
14	10.200	1398	1402	1406	rBV	588065	880555	2.17%	0.288%
15	10.335	1416	1425	1438	rBV	906504	2322333	5.72%	0.759%
16	10.494	1438	1452	1458	rBV	3218308	5891393	14.52%	1.925%
17	10.565	1458	1464	1470	rBV	736268	1372633	3.38%	0.448%
18	10.741	1487	1494	1501	rVB	864739	1575454	3.88%	0.515%
19	10.924	1517	1525	1530	rBV5	157102	406494	1.00%	0.133%
20	11.006	1530	1539	1547	rVV2	650999	1499956	3.70%	0.490%
21	11.082	1547	1552	1558	rVB2	266249	479995	1.18%	0.157%
22	11.171	1558	1567	1574	rBV	329558	666308	1.64%	0.218%
23	11.271	1574	1584	1592	rBV	528984	1137977	2.81%	0.372%
24	11.571	1631	1635	1639	rBV	274601	458885	1.13%	0.150%
25	11.641	1639	1647	1659	rVB2	1187256	2726334	6.72%	0.891%
26	11.741	1659	1664	1674	rBV	888716	1585191	3.91%	0.518%
27	11.859	1682	1684	1689	rVB2	740399	1036257	2.55%	0.339%
28	11.929	1689	1696	1698	rVV	478184	966114	2.38%	0.316%
29	11.965	1698	1702	1708	rVB	1273422	2070515	5.10%	0.676%
30	12.065	1711	1719	1731	rBV	15763904	30652217	75.56%	10.014%
31	12.300	1753	1759	1764	rBV3	323308	537869	1.33%	0.176%
32	12.394	1766	1775	1782	rVB	2154395	4039624	9.96%	1.320%
33	12.476	1784	1789	1795	rVB2	437522	775180	1.91%	0.253%
34	12.694	1820	1826	1831	rBV	1489380	2568168	6.33%	0.839%
35	12.824	1840	1848	1855	rVV	2128146	3956495	9.75%	1.293%
36	12.876	1855	1857	1861	rVB	471883	548656	1.35%	0.179%

LSC Area Percent Report

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Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Title : SW846 8260

37	12.929	1861	1866	1871	rVB3	336094	616252	1.52%	0.201%
38	12.988	1871	1876	1879	rBV	694891	1243210	3.06%	0.406%
39	13.035	1879	1884	1888	rVB	1698028	2697055	6.65%	0.881%
40	13.094	1888	1894	1898	rBV	14820843	25133716	61.96%	8.211%
41	13.129	1898	1900	1903	rVV	6198338	8416300	20.75%	2.750%
42	13.171	1903	1907	1917	rVB	8790911	15455062	38.10%	5.049%
43	13.335	1926	1935	1942	rBV	4765306	8427727	20.78%	2.753%
44	13.423	1942	1950	1954	rBV	1732637	3131874	7.72%	1.023%
45	13.482	1954	1960	1966	rBV	25881559	40564885	100.00%	13.253%
46	13.529	1966	1968	1974	rVB	761096	1069748	2.64%	0.349%
47	13.641	1974	1987	1991	rBV3	2060374	5288215	13.04%	1.728%
48	13.718	1996	2000	2002	rBV2	336420	443281	1.09%	0.145%
49	13.823	2008	2018	2031	rVB2	6778569	16993265	41.89%	5.552%
50	13.953	2034	2040	2041	rBV3	1386840	2070712	5.10%	0.677%
51	13.988	2041	2046	2049	rBV2	4469104	8799693	21.69%	2.875%
52	14.106	2062	2066	2074	rBV2	647339	1167136	2.88%	0.381%
53	14.182	2074	2079	2084	rVB	1612165	2632449	6.49%	0.860%
54	14.241	2084	2089	2092	rBV	2491024	4228645	10.42%	1.382%
55	14.270	2092	2094	2099	rVV	2343977	3093239	7.63%	1.011%
56	14.329	2099	2104	2112	rVB	4386174	7036700	17.35%	2.299%
57	14.441	2117	2123	2128	rVB3	1150033	2017032	4.97%	0.659%
58	14.570	2140	2145	2151	rBV	969560	1564541	3.86%	0.511%
59	14.653	2151	2159	2162	rBV	2923524	5055306	12.46%	1.652%
60	14.694	2162	2166	2170	rVB	3183129	4425207	10.91%	1.446%
61	14.735	2170	2173	2177	rVB2	1266248	1626533	4.01%	0.531%
62	14.806	2182	2185	2192	rVB	646876	948478	2.34%	0.310%
63	14.876	2192	2197	2200	rBV	574557	787649	1.94%	0.257%
64	14.923	2200	2205	2209	rBV	1361251	2527481	6.23%	0.826%
65	14.959	2209	2211	2216	rVB2	1086752	1428797	3.52%	0.467%
66	15.029	2216	2223	2224	rBV3	2222778	4060989	10.01%	1.327%
67	15.094	2232	2234	2243	rVB3	585521	1037164	2.56%	0.339%
68	15.212	2249	2254	2260	rVB4	245019	520788	1.28%	0.170%
69	15.317	2266	2272	2284	rVB3	985685	2519175	6.21%	0.823%
70	15.429	2285	2291	2300	rVB4	652529	1640225	4.04%	0.536%
71	15.594	2313	2319	2320	rBV3	335122	511113	1.26%	0.167%
72	15.635	2321	2326	2334	rVB	2057010	4002148	9.87%	1.308%
73	15.747	2339	2345	2349	rBV2	343940	677301	1.67%	0.221%
74	15.829	2350	2359	2369	rVB3	604028	1774023	4.37%	0.580%
75	15.941	2370	2378	2386	rBV	524936	1287177	3.17%	0.421%
76	16.123	2405	2409	2421	rVB3	249673	745408	1.84%	0.244%

LSC Area Percent Report

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Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

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\82N091024W.M
Title : SW846 8260

77	16.253	2426	2431	2437	rVV3	267135	596564	1.47%	0.195%
78	16.335	2438	2445	2452	rVB3	359988	910548	2.24%	0.297%
79	16.782	2513	2521	2522	rBV	1394668	2490563	6.14%	0.814%

Sum of corrected areas: 306081708

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
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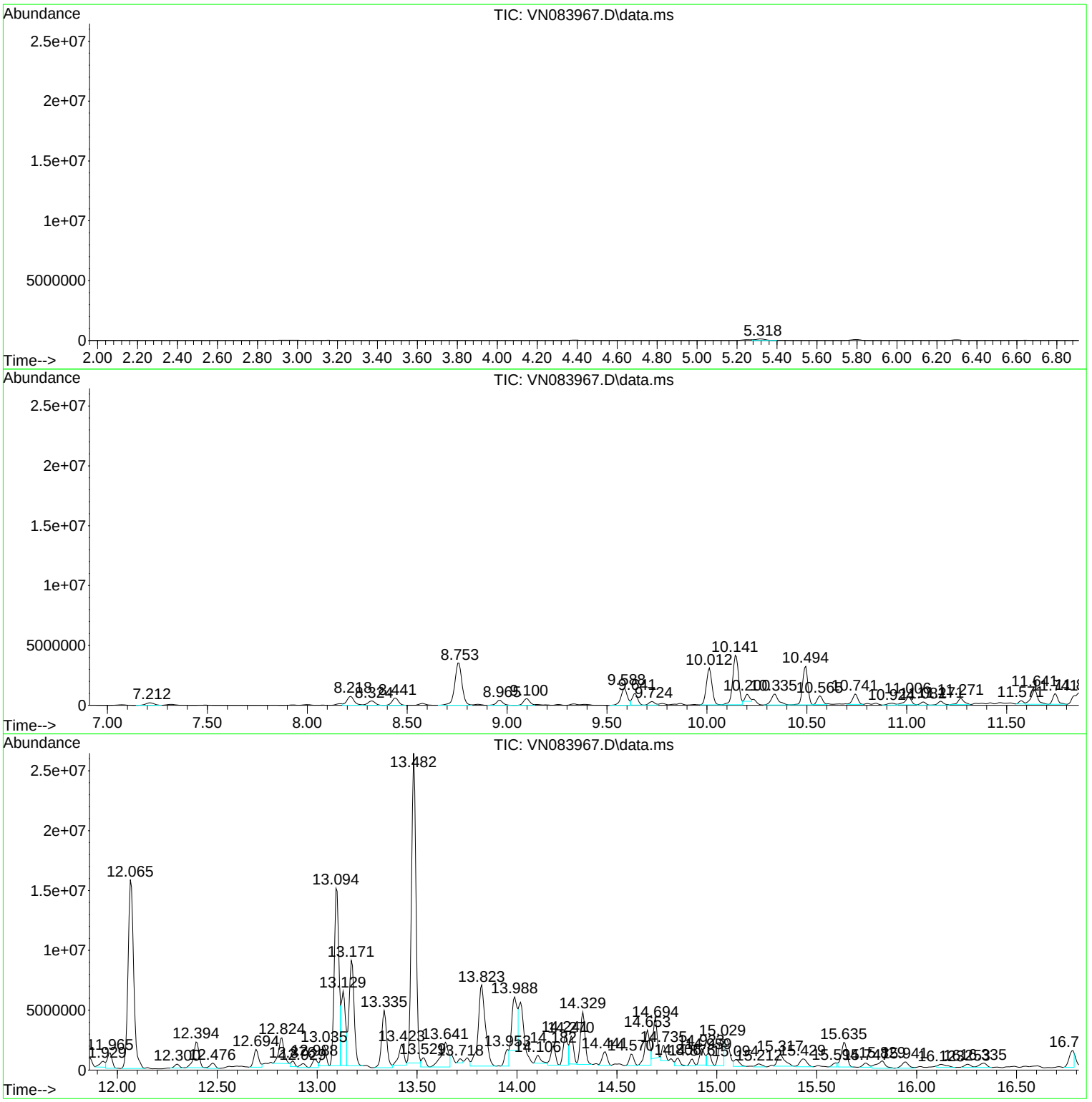
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

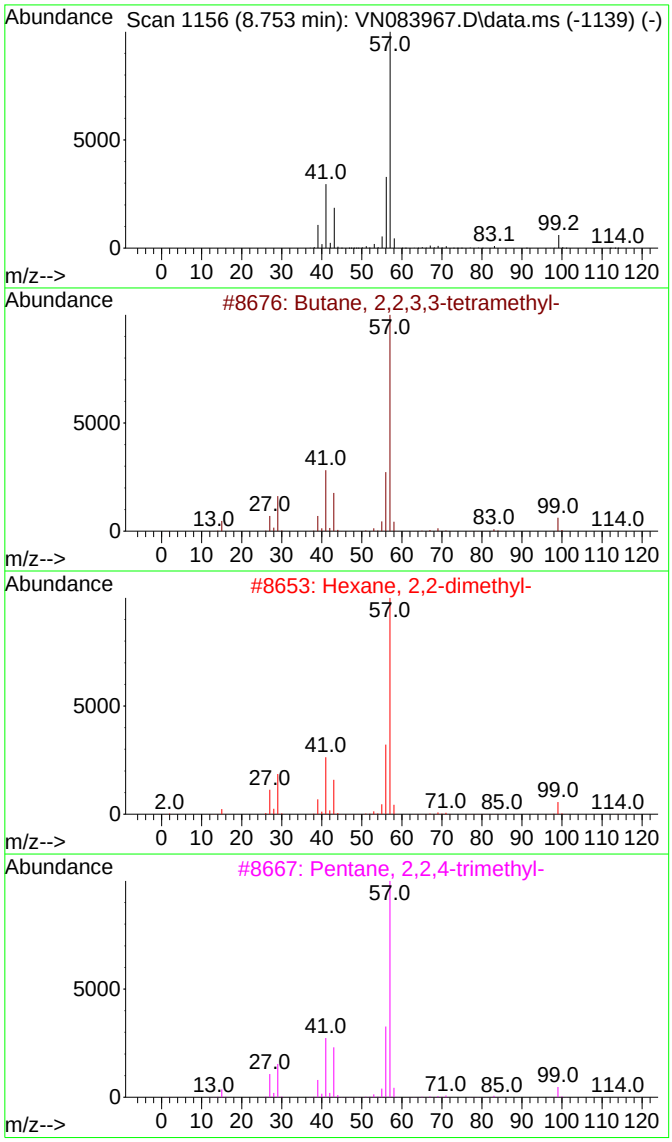
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

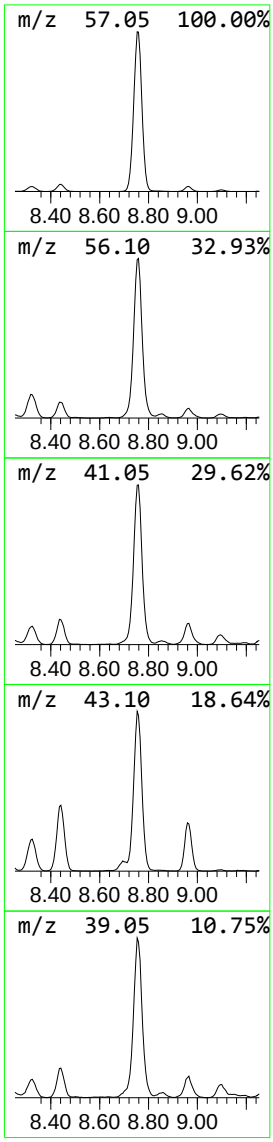
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.753	355.29 ug/l	8874180	1,4-Difluorobenzene	9.100

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	78
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	72
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

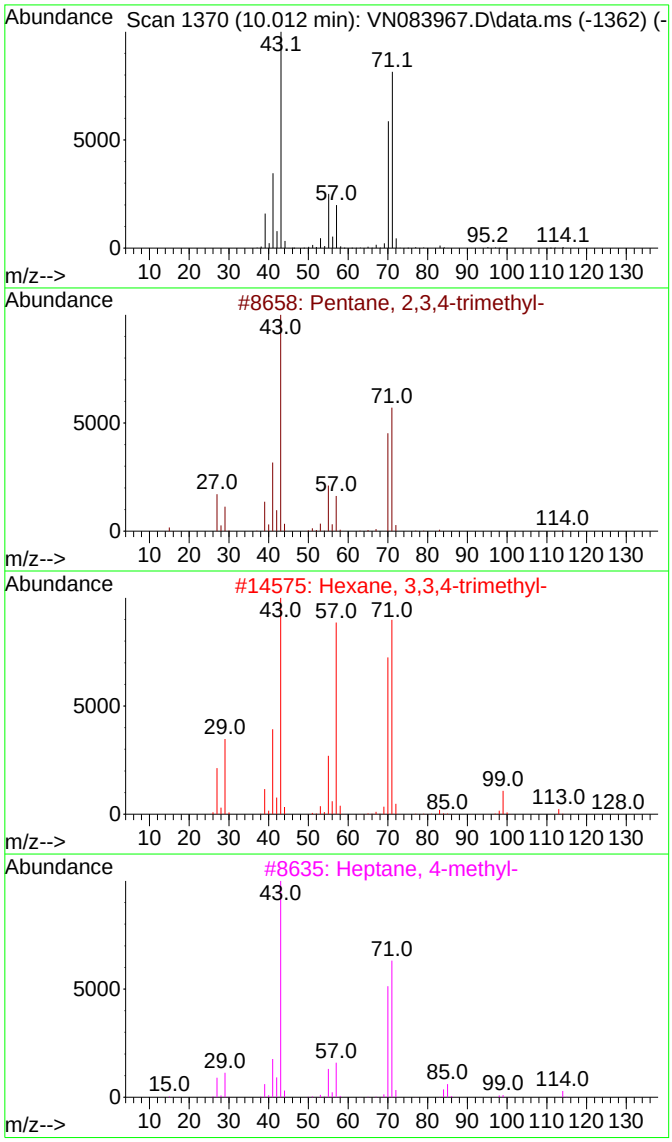
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

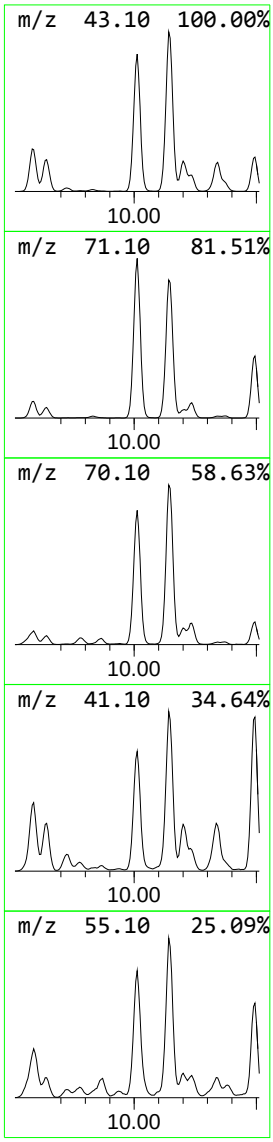
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.012	241.24 ug/l	6025550	1,4-Difluorobenzene	9.100

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	83
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
4	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

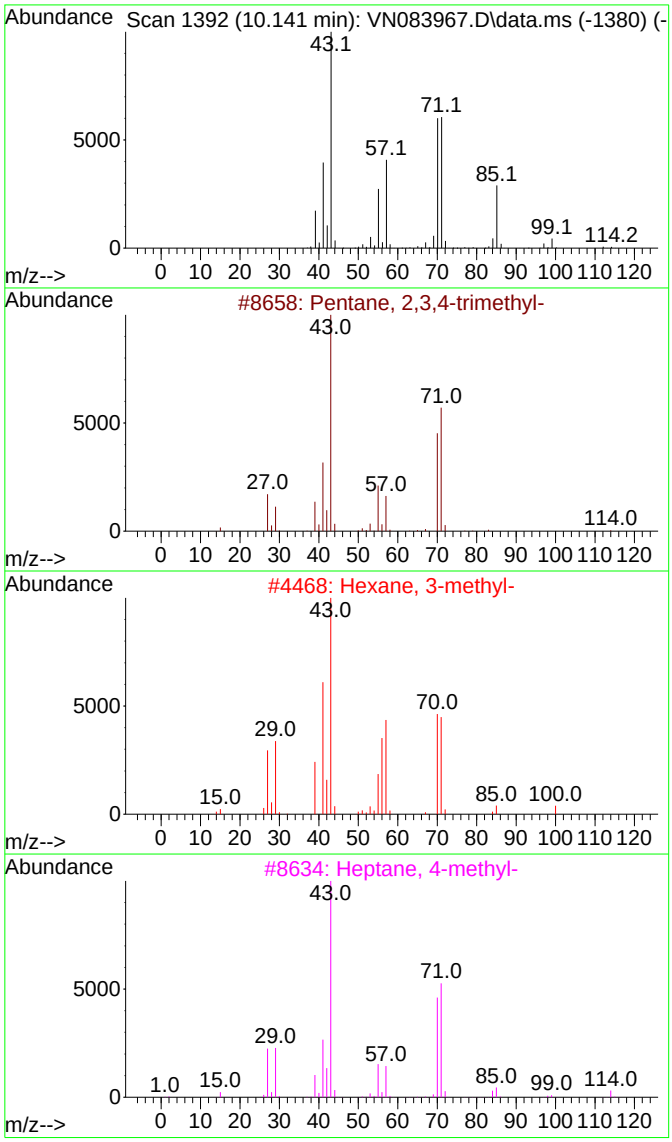
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

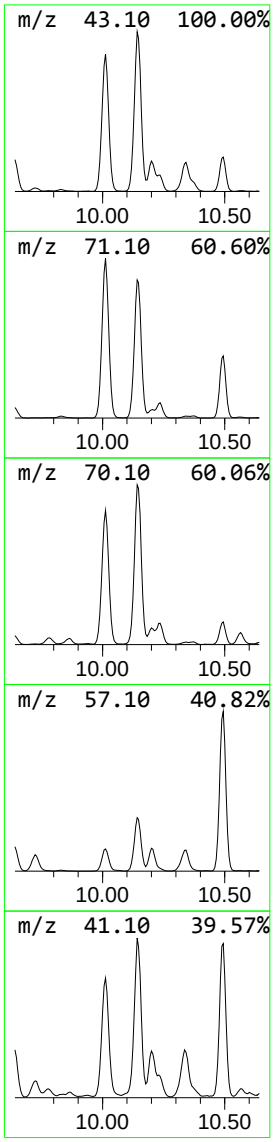
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.141	336.32 ug/l	8400240	1,4-Difluorobenzene	9.100

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	81
2	Hexane, 3-methyl-	100	C7H16	000589-34-4	64
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	64
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
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Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

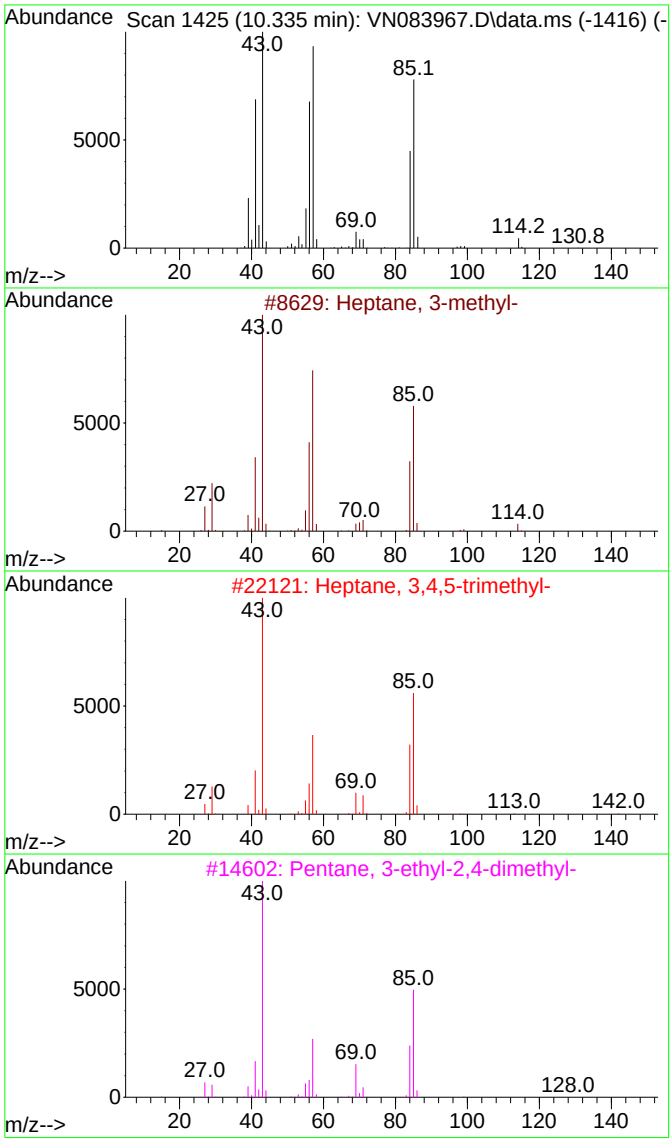
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

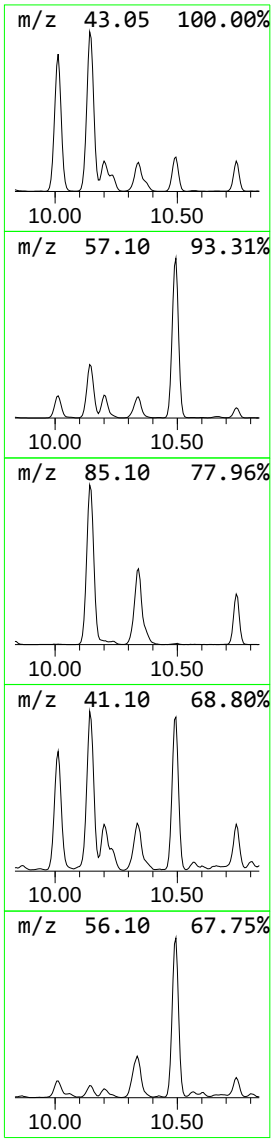
****

Peak Number 4 Heptane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.335	92.98 ug/l	2322330	1,4-Difluorobenzene	9.100

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	64
3	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59
4	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	58
5	Heptane, 4-ethyl-	128	C9H20	002216-32-2	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

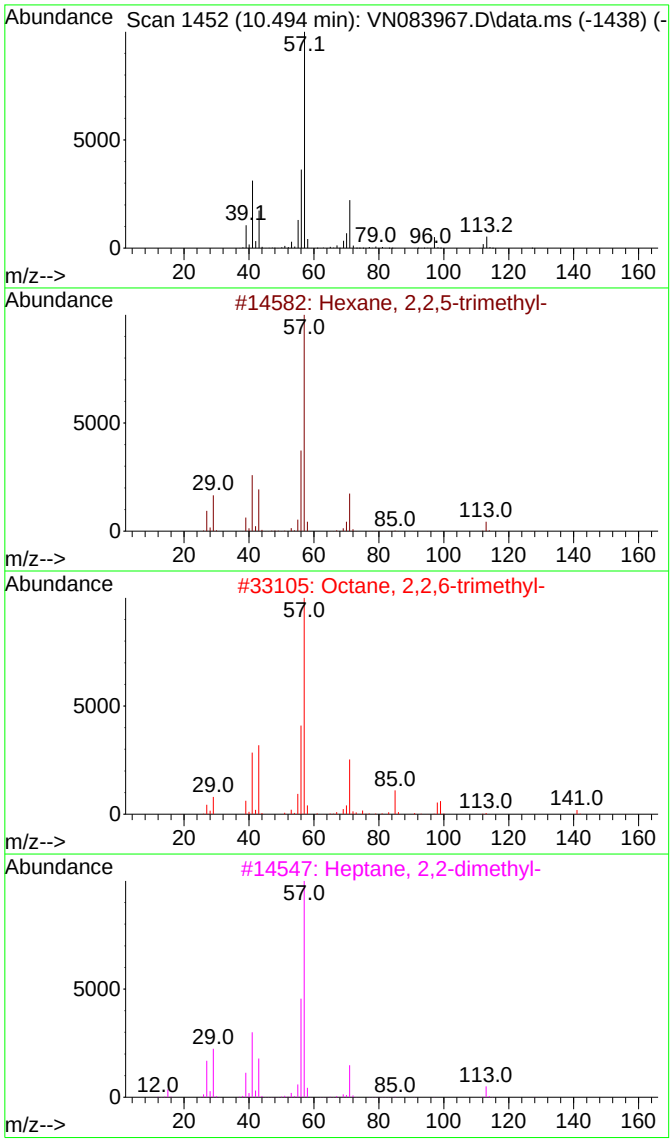
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

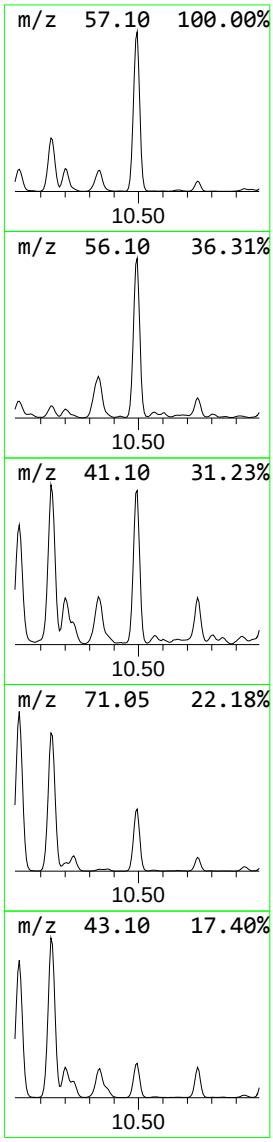
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexane, 2,2,5-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.494	284.26 ug/l	5891390	Chlorobenzene-d5	11.865

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	72
2	Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	64
3	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	53
4	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	45
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

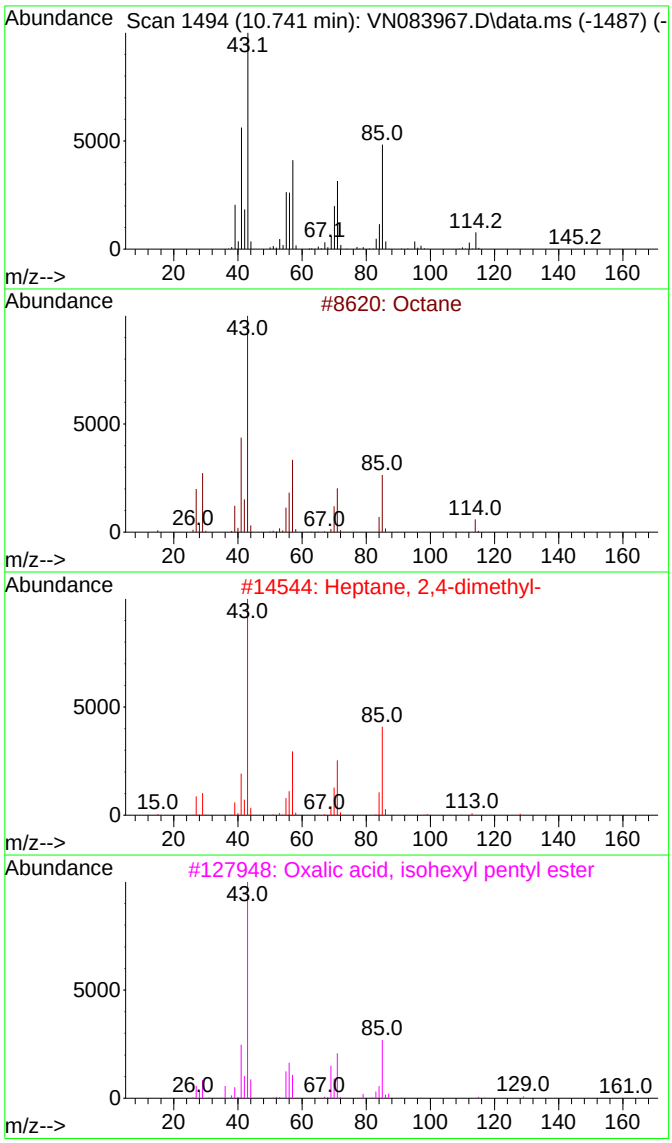
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

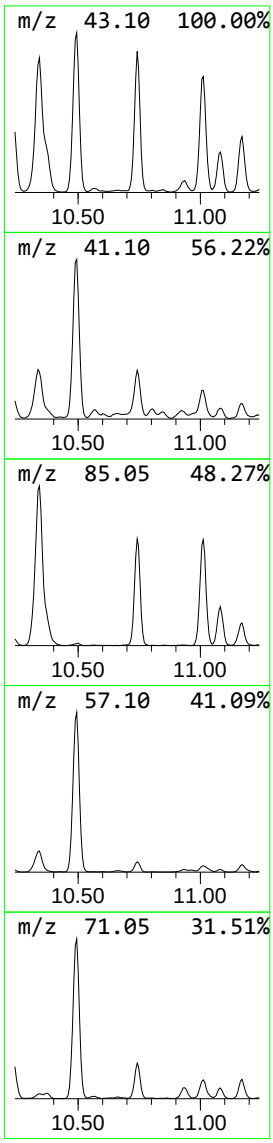
****

Peak Number 6 Octane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.741	76.02 ug/l	1575450	Chlorobenzene-d5	11.865

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	91
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	80
3	Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	72
4	Undecane, 2,4-dimethyl-	184	C13H28	017312-80-0	59
5	Hexane, 3-ethyl-	114	C8H18	000619-99-8	59





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Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

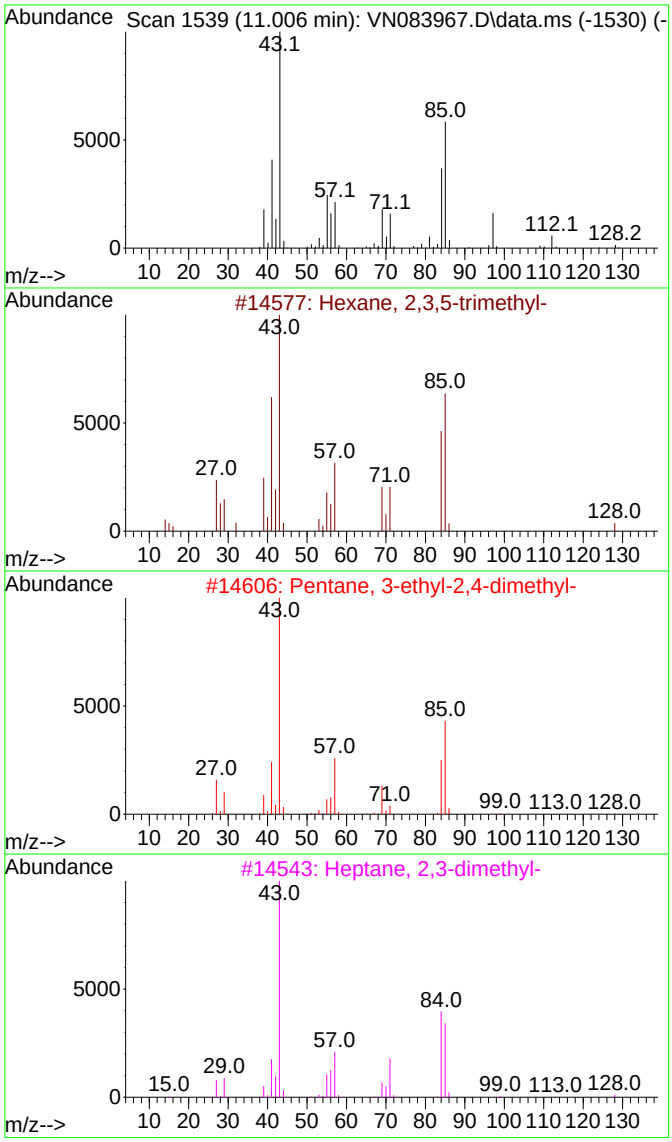
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

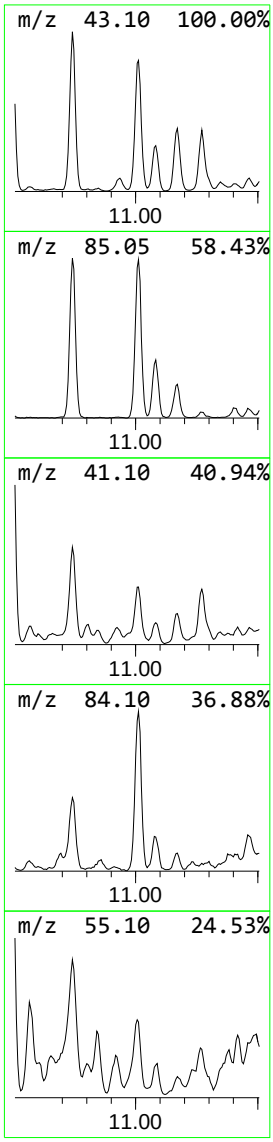
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexane, 2,3,5-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.006	72.37 ug/l	1499960	Chlorobenzene-d5	11.865

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	72
2	Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	64
3	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	62
4	Nonane, 5-methyl-	142	C10H22	015869-85-9	59
5	Pentane, 2,3,3,4-tetramethyl-	128	C9H20	016747-38-9	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

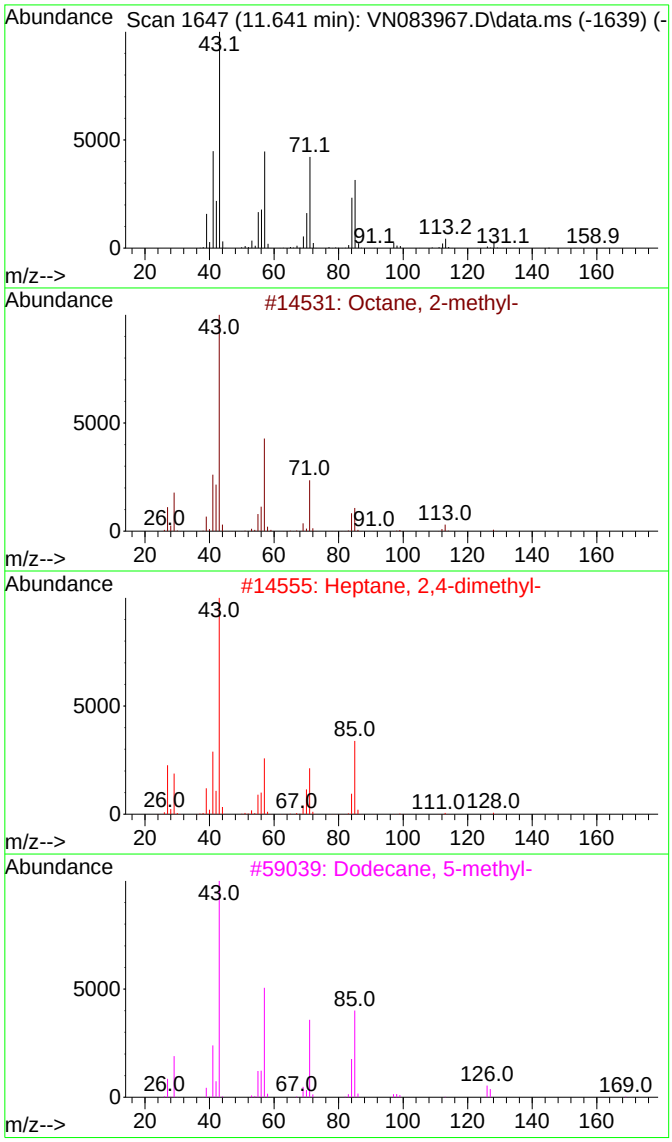
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

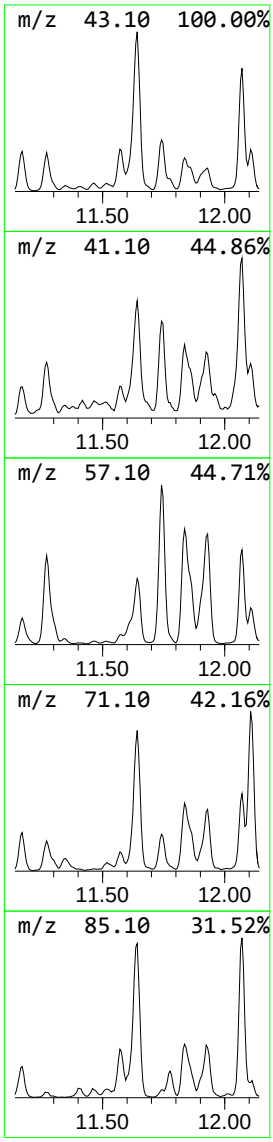
****

Peak Number 8 Octane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.641	131.55 ug/l	2726330	Chlorobenzene-d5	11.865

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	80
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3	Dodecane, 5-methyl-	184	C13H28	017453-93-9	59
4	Undecane, 5-methyl-	170	C12H26	001632-70-8	59
5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

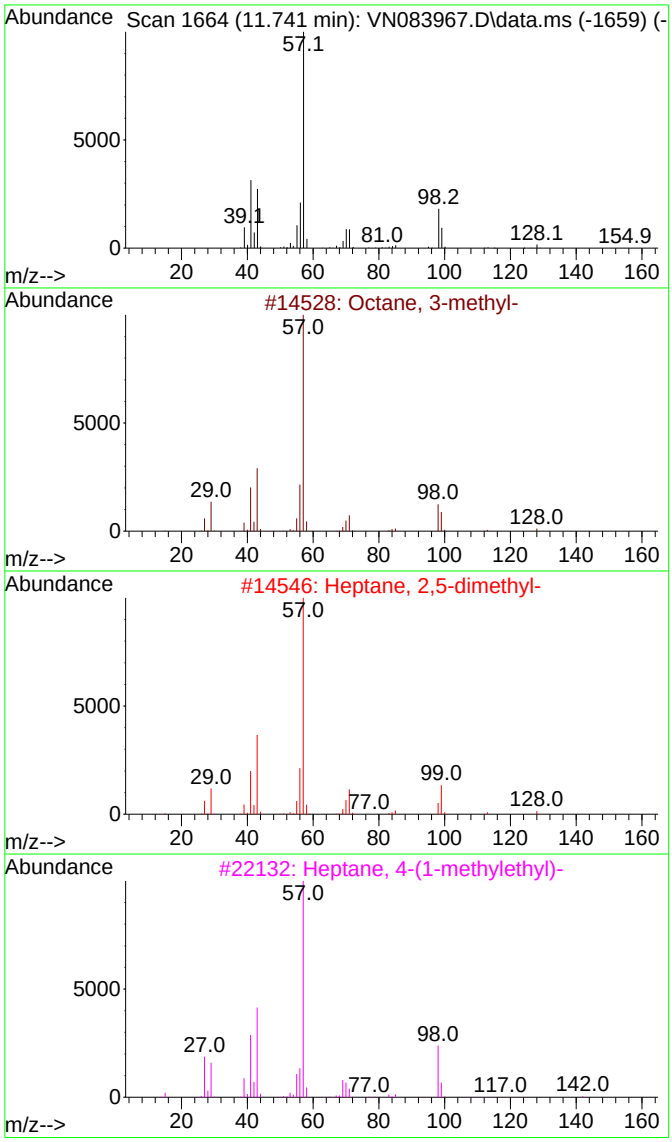
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

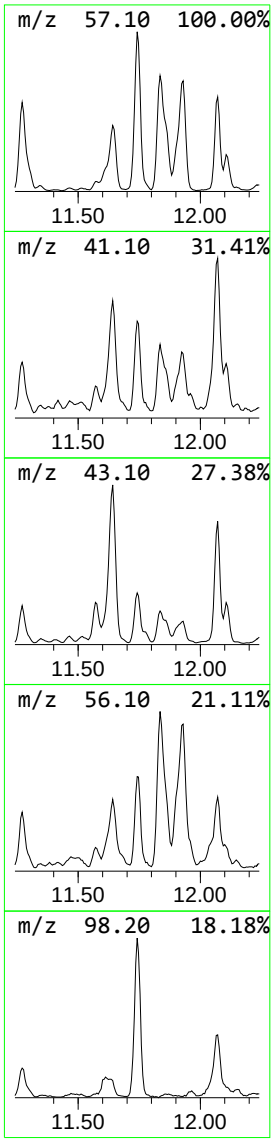
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Octane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.741	76.49 ug/l	1585190	Chlorobenzene-d5	11.865

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-	128	C9H20	002216-33-3	91
2	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	87
3	Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	72
4	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	52
5	Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

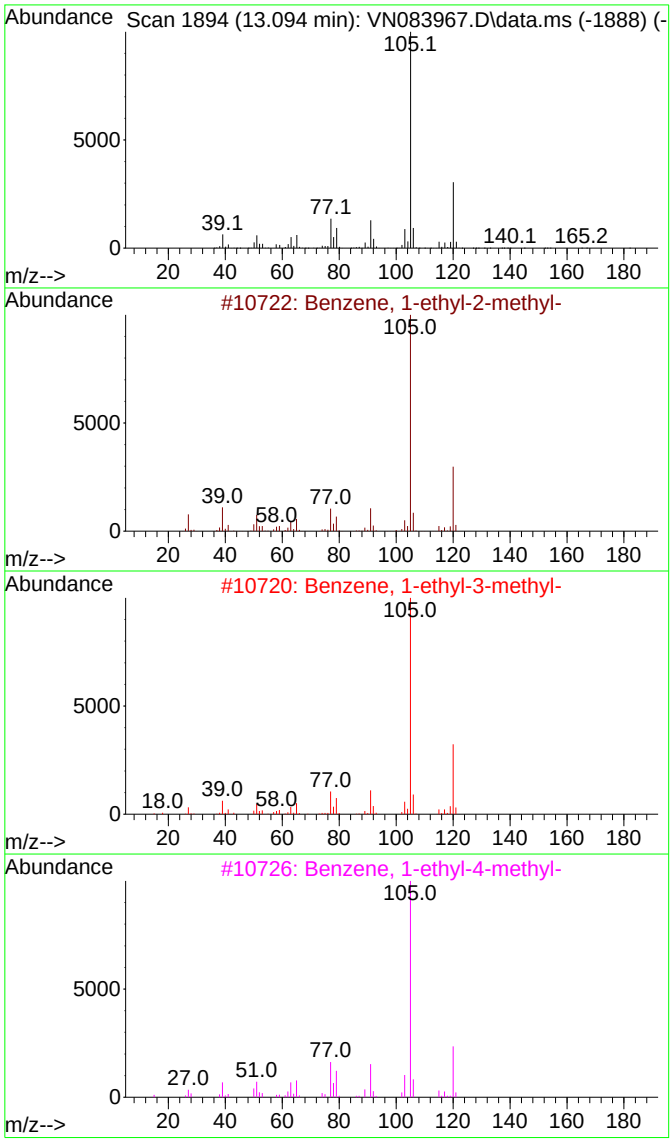
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
Quant Title : SW846 8260

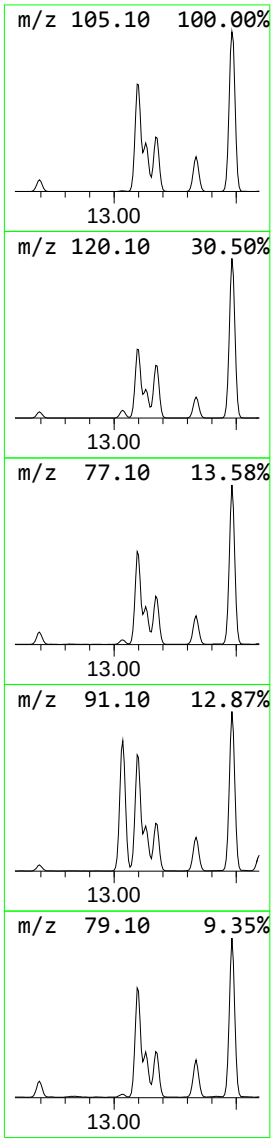
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.094	73.95 ug/l	25133700	1,4-Dichlorobenzene-d4	13.788

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	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
Data File : VN083967.D
Acq On : 18 Sep 2024 07:09
Operator : JC\MD
Sample : P3964-06
Misc : 6.26g/10mL/100uL/5.00ml/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SINK-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.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
Butane, 2,2,3,3...	8.753	355.3	ug/l	8874180	2	9.100	1248850	50.0
Pentane, 2,3,4-...	10.012	241.2	ug/l	6025550	2	9.100	1248850	50.0
Hexane, 3-methyl-	10.141	336.3	ug/l	8400240	2	9.100	1248850	50.0
Heptane, 3-methyl-	10.335	93.0	ug/l	2322330	2	9.100	1248850	50.0
Hexane, 2,2,5-t...	10.494	284.3	ug/l	5891390	3	11.865	1036260	50.0
Octane	10.741	76.0	ug/l	1575450	3	11.865	1036260	50.0
Hexane, 2,3,5-t...	11.006	72.4	ug/l	1499960	3	11.865	1036260	50.0
Octane, 2-methyl-	11.641	131.6	ug/l	2726330	3	11.865	1036260	50.0
Octane, 3-methyl-	11.741	76.5	ug/l	1585190	3	11.865	1036260	50.0
Benzene, 1-ethy...	13.094	74.0	ug/l	25133700	4	13.788	16993300	50.0