LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

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\82Y081423S.M
Title : SW846 8260

Signal : TIC: VY015196.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.071	35	41	61	rBV	11666482	18256664	6.43%	0.360%
2	2.235	61	68	99	rVV	15749122	46220373	16.27%	0.911%
3	2.875	165	173	211	rBV2	25920390	137111286	48.27%	2.702%
4	3.156	211	219	223	rBV	2055384	4298680	1.51%	0.085%
5	3.430	256	264	279	rVB	5936576	14583643	5.13%	0.287%
6	3.595	279	291	298	rVV	2665882	7087600	2.50%	0.140%
7	3.680	298	305	332	rVB	8813299	24272481	8.55%	0.478%
8	3.930	332	346	386	rVB	9785484	34007893	11.97%	0.670%
9	4.576	430	452	459	rBV6	1453131	6115715	2.15%	0.121%
10	4.777	459	485	537	rVB6	30134544	284042381	100.00%	5.598%
11	5.228	542	559	561	rBV	25768473	72696180	25.59%	1.433%
12	5.570	599	615	631	rBV4	3786990	17418902	6.13%	0.343%
13	5.741	631	643	644	rBV2	25727977	55492201	19.54%	1.094%
14	5.917	664	672	681	rBV	3072600	9070657	3.19%	0.179%
15	6.039	681	692	697	rVV	4598773	15575649	5.48%	0.307%
16	6.100	697	702	715	rVB	6117110	19736038	6.95%	0.389%
17	6.240	715	725	734	rBV	3224593	11223957	3.95%	0.221%
18	6.375	735	747	762	rVB2	2307896	10516365	3.70%	0.207%
19	6.545	762	775	788	rBV2	10459500	39317363	13.84%	0.775%
20	6.704	788	801	809	rBV	18705899	75111961	26.44%	1.480%
21	6.801	809	817	839	rVB2	26915048	110816537	39.01%	2.184%
22	6.990	839	848	879	rVB	1372878	5515108	1.94%	0.109%
23	7.539	918	938	949	rBV3	10848234	42190323	14.85%	0.831%
24	7.795	966	980	989	rBV4	29295803	153871869	54.17%	3.032%
25	7.899	989	997	1007	rVB2	27117019	103841615	36.56%	2.046%
26	8.033	1007	1019	1036	rVB2	31951201	150039163	52.82%	2.957%
27	8.362	1036	1073	1081	rBV6	29225767	209123490	73.62%	4.121%
28	8.435	1081	1085	1096	rVB	12958142	33679685	11.86%	0.664%
29	8.582	1096	1109	1120	rVB2	31061935	127236122	44.79%	2.507%
30	8.703	1120	1129	1134	rBV2	14759395	38681394	13.62%	0.762%
31	8.862	1149	1155	1162	rVB	3695827	7965329	2.80%	0.157%
32	8.972	1162	1173	1176	rBV6	3968561	13045126	4.59%	0.257%
33	9.014	1177	1180	1194	rVB	5194091	13608408	4.79%	0.268%
34	9.197	1194	1210	1216	rBV9	32264381	133652210	47.05%	2.634%
35	9.264	1217	1221	1229	rVV	26732818	70160686	24.70%	1.383%
36	9.344	1229	1234	1236	rVV	5197605	10757815	3.79%	0.212%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

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\82Y081423S.M
Title : SW846 8260

37	9.380	1236	1240	1246	rVV2	8517175	19670311	6.93%	0.388%
38	9.466	1246	1254	1266	rVB4	8413085	29500413	10.39%	0.581%
39	9.636	1272	1282	1292	rBV3	28630674	108257811	38.11%	2.133%
40	9.770	1292	1304	1311	rBV7	29500128	128139334	45.11%	2.525%
41	9.862	1311	1319	1330	rVB5	26696362	131356275	46.25%	2.589%
42	9.996	1330	1341	1354	rVB7	28126998	136659204	48.11%	2.693%
43	10.173	1358	1370	1376	rBV	16958877	51925838	18.28%	1.023%
44	10.258	1376	1384	1389	rBV6	31328167	114762452	40.40%	2.262%
45	10.392	1399	1406	1413	rVB2	29896775	86239559	30.36%	1.700%
46	10.447	1413	1415	1420	rVB	2919142	4102137	1.44%	0.081%
47	10.545	1427	1431	1438	rVB3	2251345	5805976	2.04%	0.114%
48	10.642	1438	1447	1456	rBV7	11255296	40165703	14.14%	0.792%
49	10.728	1456	1461	1468	rVB	11462357	24713549	8.70%	0.487%
50	10.819	1468	1476	1486	rBV2	11489405	32227026	11.35%	0.635%
51	10.923	1486	1493	1501	rBV	18002034	50386369	17.74%	0.993%
52	11.105	1518	1523	1528	rBV	4514383	10063226	3.54%	0.198%
53	11.221	1536	1542	1543	rBV	7983822	12397318	4.36%	0.244%
54	11.301	1547	1555	1565	rVB6	27114179	121603978	42.81%	2.396%
55	11.416	1565	1574	1584	rVB3	24017945	90745352	31.95%	1.788%
56	11.526	1585	1592	1597	rBV3	3626138	10137372	3.57%	0.200%
57	11.624	1597	1608	1616	rBV3	39919947	181151691	63.78%	3.570%
58	11.709	1616	1622	1628	rBV4	41235534	129008235	45.42%	2.542%
59	11.910	1650	1655	1657	rBV	2293375	3684519	1.30%	0.073%
60	11.959	1657	1663	1668	rBV3	9124895	17472270	6.15%	0.344%
61	12.044	1668	1677	1680	rBV4	42847714	117086821	41.22%	2.307%
62	12.136	1687	1692	1698	rVB	11916650	23516702	8.28%	0.463%
63	12.215	1698	1705	1708	rBV2	8571971	19298601	6.79%	0.380%
64	12.343	1720	1726	1732	rBV2	31092521	79222810	27.89%	1.561%
65	12.556	1753	1761	1766	rVB	13857420	34238532	12.05%	0.675%
66	12.691	1774	1783	1788	rBV2	37805973	129572636	45.62%	2.554%
67	12.745	1788	1792	1796	rBV2	22230672	49511335	17.43%	0.976%
68	12.989	1825	1832	1841	rVV2	48231222	129978140	45.76%	2.562%
69	13.062	1841	1844	1846	rVV2	3130159	4951172	1.74%	0.098%
70	13.087	1846	1848	1849	rVV	3757385	3825424	1.35%	0.075%
71	13.130	1849	1855	1859	rVV2	48251216	120246067	42.33%	2.370%
72	13.172	1860	1862	1868	rVV3	47783030	54274151	19.11%	1.070%
73	13.258	1868	1876	1882	rVV3	15985548	43351518	15.26%	0.854%
74	13.325	1882	1887	1891	rVV	19688070	32488884	11.44%	0.640%
75	13.379	1891	1896	1902	rVV3	7236482	14753946	5.19%	0.291%
76	13.471	1902	1911	1925	rVB2	44979763	111047478	39.10%	2.188%

LSC Area Percent Report

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Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

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\82Y081423S.M
Title : SW846 8260

77	13.599	1925	1932	1933	rBV	20547141	29700118	10.46%	0.585%
78	13.629	1933	1937	1941	rVV3	45260423	98870481	34.81%	1.948%
79	13.672	1941	1944	1954	rVB4	38242852	86895683	30.59%	1.712%
80	13.757	1954	1958	1963	rBV2	5969727	8404413	2.96%	0.166%
81	13.825	1963	1969	1974	rVB	16970354	26500368	9.33%	0.522%
82	13.885	1974	1979	1982	rBV	20611058	34183734	12.03%	0.674%
83	13.916	1982	1984	1989	rVV	18077483	22875999	8.05%	0.451%
84	13.971	1989	1993	1999	rVB	26888116	41063356	14.46%	0.809%
85	14.081	2006	2011	2016	rVB2	10118214	16154692	5.69%	0.318%
86	14.215	2027	2033	2037	rBV	5446669	8002831	2.82%	0.158%
87	14.294	2037	2046	2049	rBV	8915158	14557129	5.12%	0.287%
88	14.337	2049	2053	2057	rVV	11123756	16189938	5.70%	0.319%
89	14.379	2057	2060	2063	rVV	3383331	4409357	1.55%	0.087%
90	14.562	2086	2090	2095	rVV2	4223255	6895155	2.43%	0.136%
91	14.611	2095	2098	2102	rVB	2325679	3106845	1.09%	0.061%
92	14.678	2102	2109	2115	rBV3	5738657	13368850	4.71%	0.263%
93	14.940	2147	2152	2164	rVB2	1390880	3356983	1.18%	0.066%
94	15.233	2189	2200	2206	rBV	3343503	5874658	2.07%	0.116%

Sum of corrected areas: 5074291594

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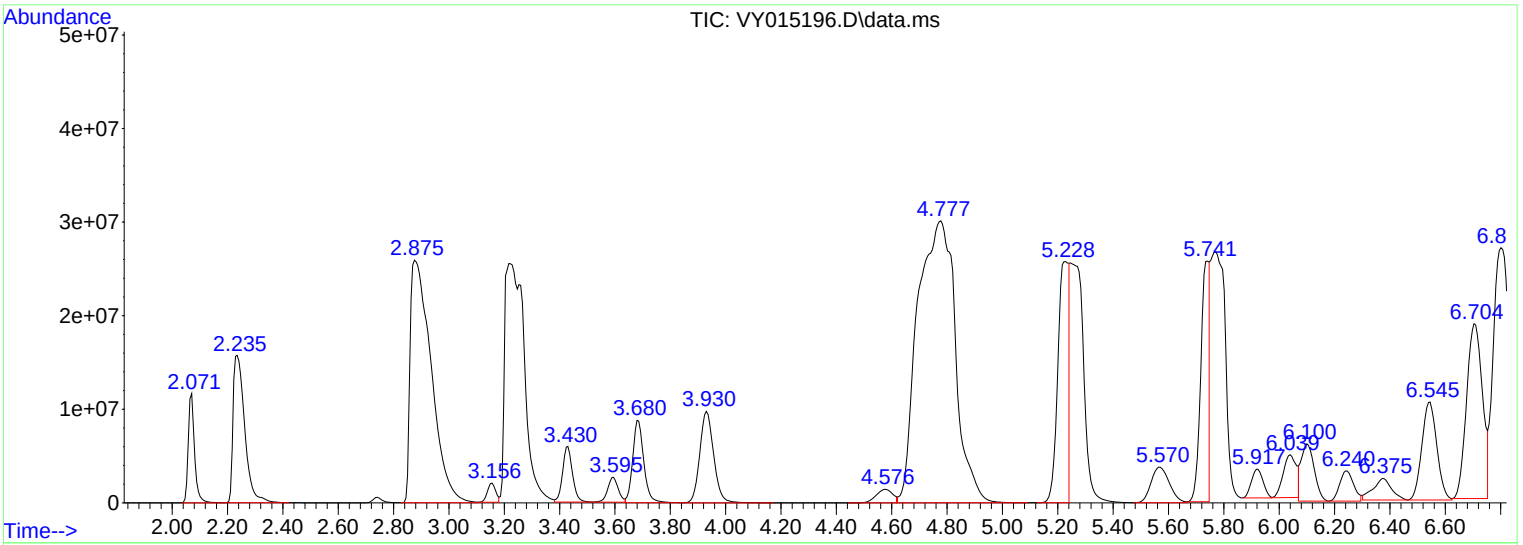
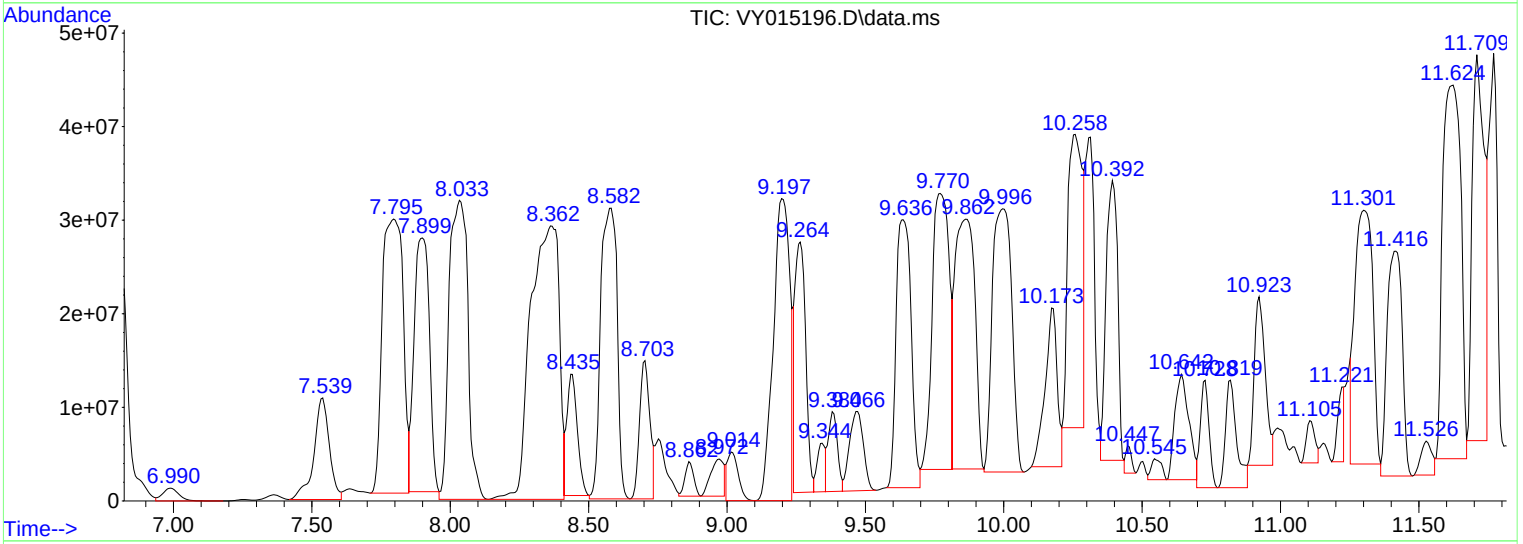
ALS Vial : 15 Sample Multiplier: 1

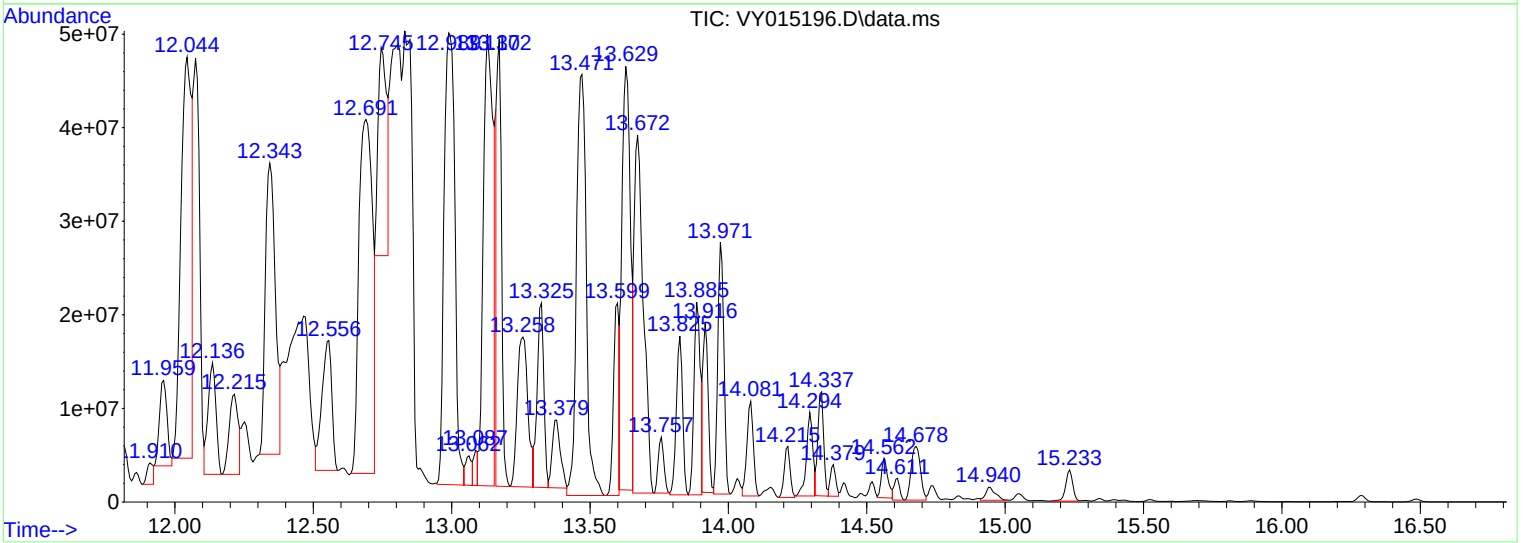
Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

Quant Title :

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







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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

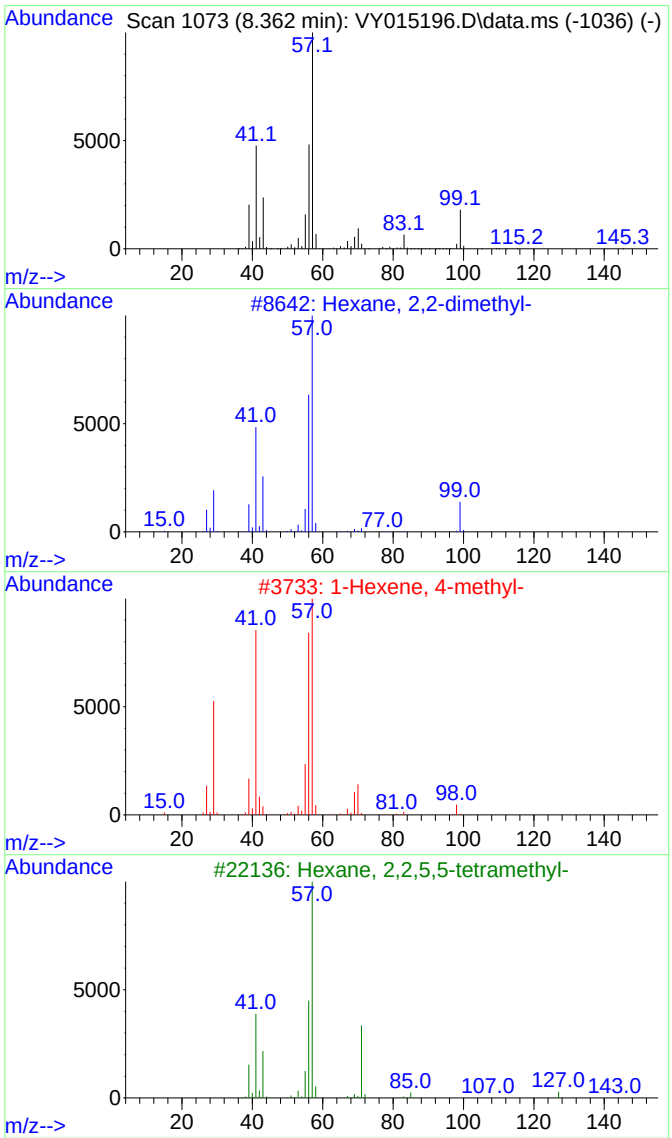
Quant Title :

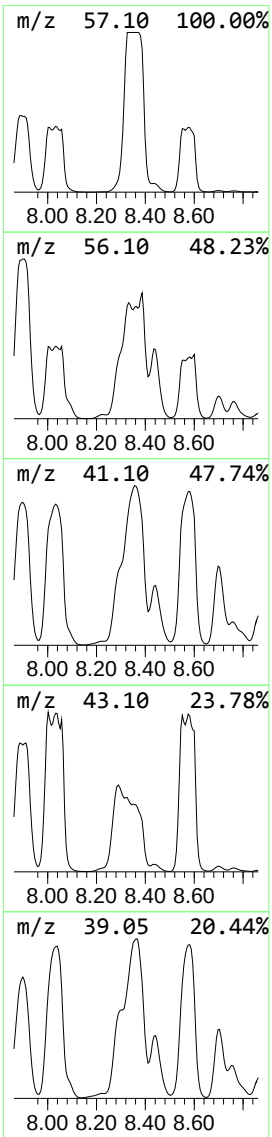
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Hexane, 2,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.362	270.32 ug/l	209123000	1,4-Difluorobenzene	8.697

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	53
2	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	53
3	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	50
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	45
5	Butane, 2-azido-2,3,3-trimethyl-	141	C7H15N3	051677-41-9	42





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Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

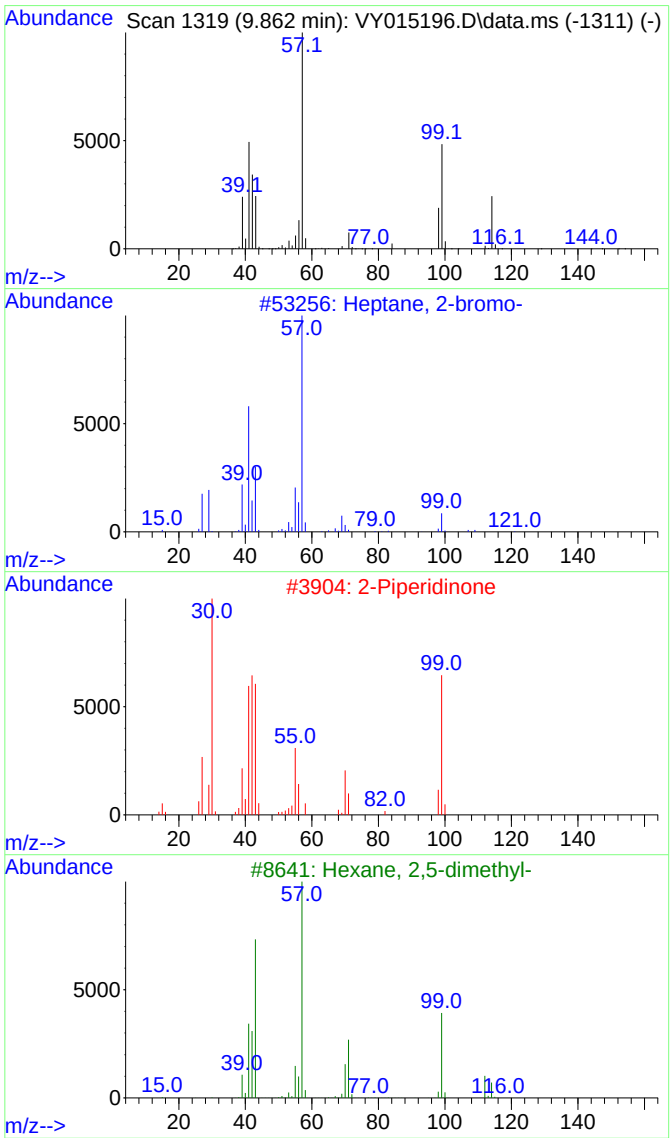
Quant Title :

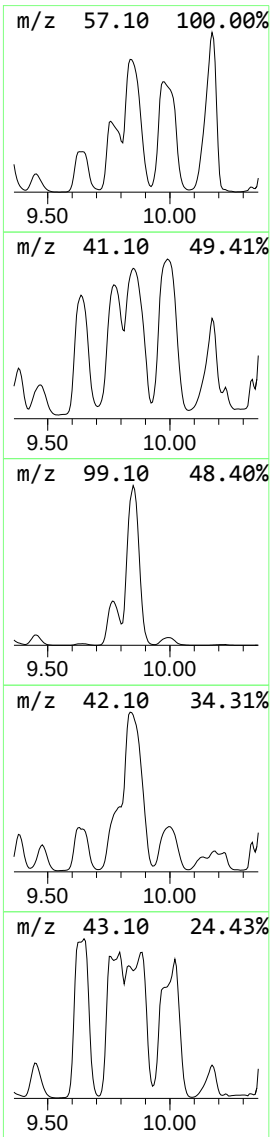
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 unknown9.862 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.862	169.79 ug/l	131356000	1,4-Difluorobenzene	8.697

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-bromo-	178	C7H15Br	001974-04-5	47
2	2-Piperidinone	99	C5H9NO	000675-20-7	47
3	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	45
4	Acetamide, N-ethenyl-N-methyl-	99	C5H9NO	003195-78-6	38
5	Heptane, 3-bromo-	178	C7H15Br	001974-05-6	37





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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

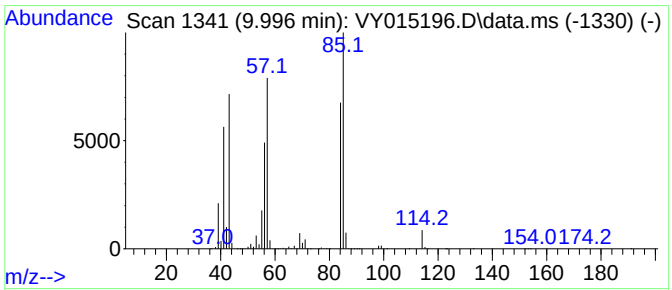
Quant Title :

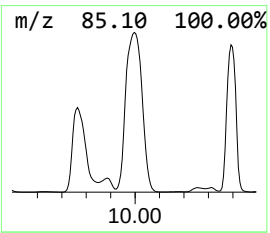
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

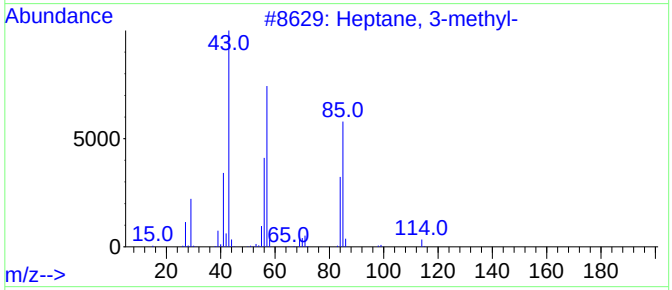
****
Peak Number 3 Heptane, 3-methyl- Concentration Rank 9

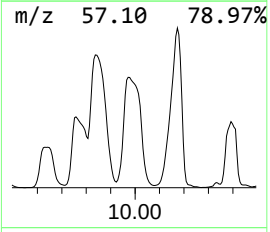
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.996	176.65 ug/l	136659000	1,4-Difluorobenzene	8.697

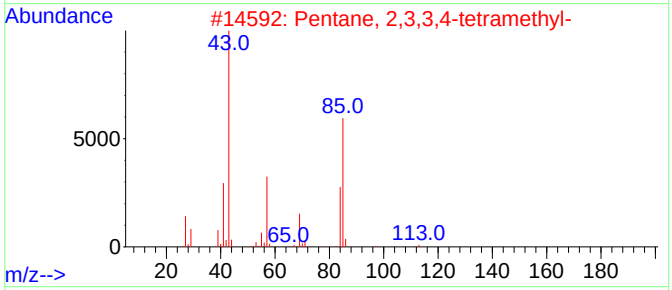
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	78
2	Pentane, 2,3,3,4-tetramethyl-	128	C9H20	016747-38-9	59
3	Nonane, 5-methyl-	142	C10H22	015869-85-9	59
4	Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	59
5	Hexane, 2-methyl-	100	C7H16	000591-76-4	59

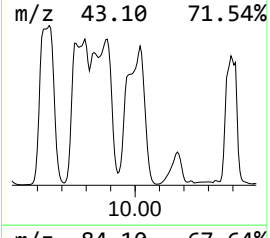


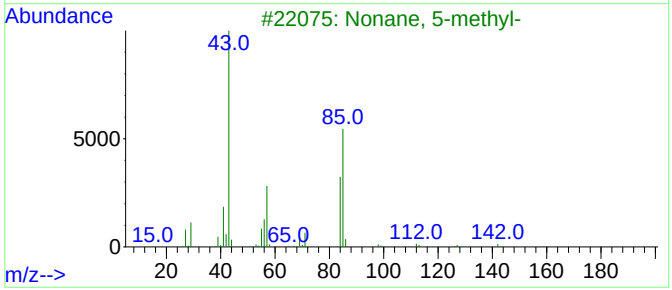


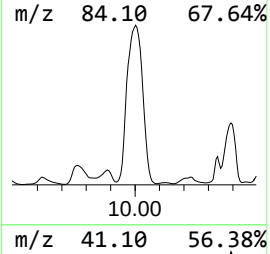


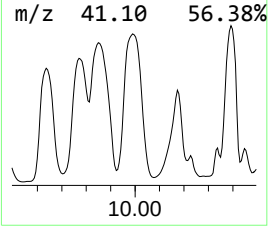












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ClientSampleId :
B-PP09C

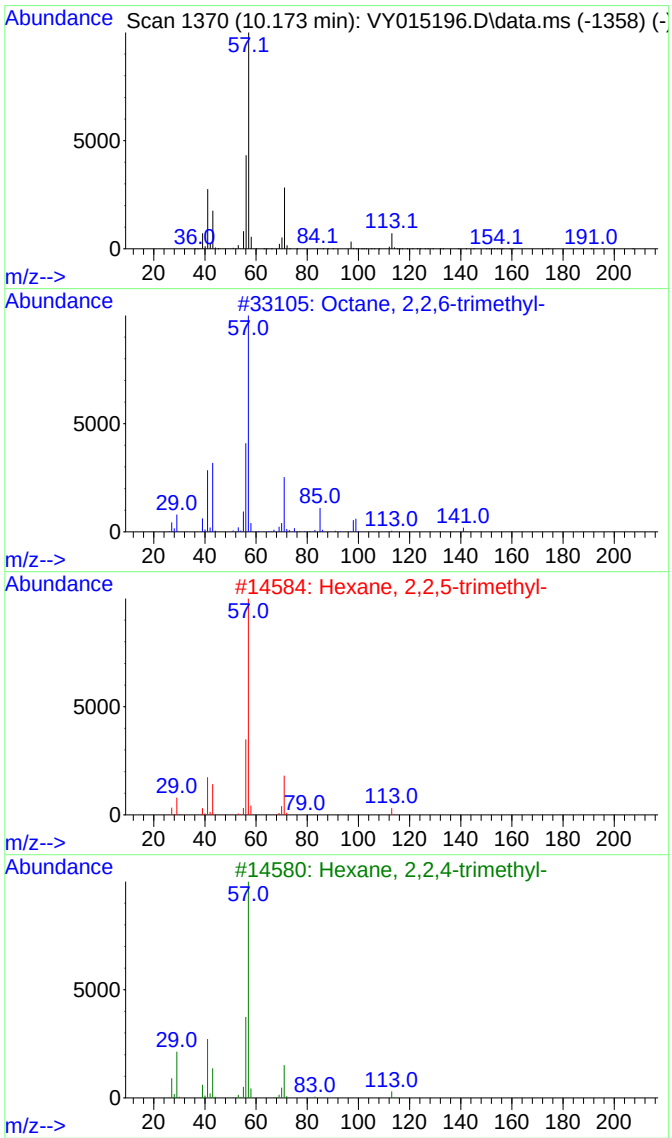
Quant Title :

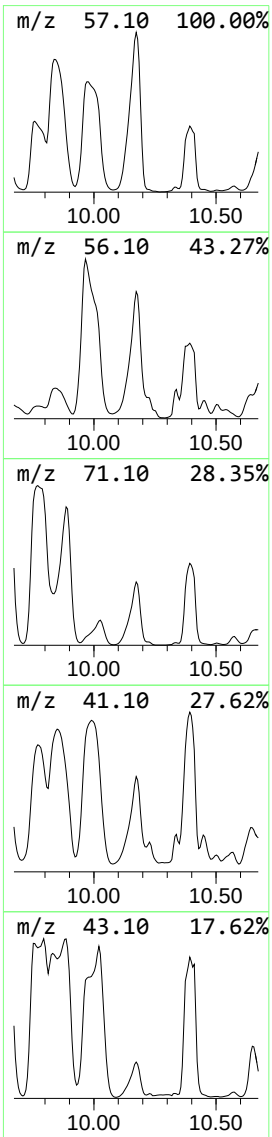
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Octane, 2,2,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.173	256.11 ug/l	51925800	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	78
2	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	78
3	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	72
4	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	72
5	Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

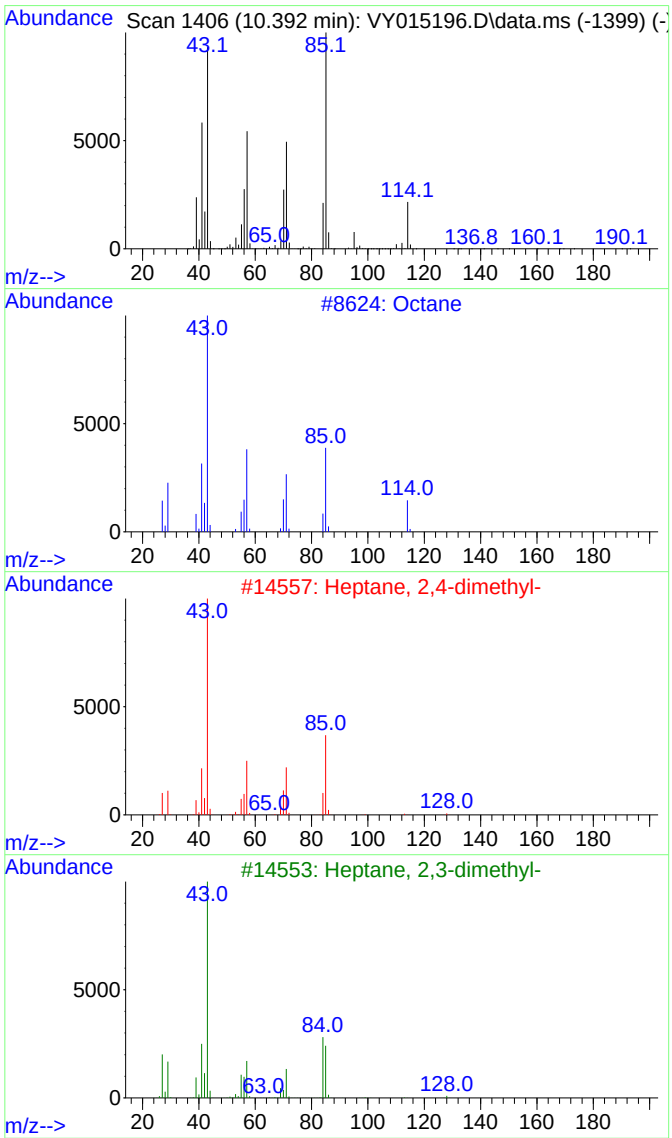
Quant Title :

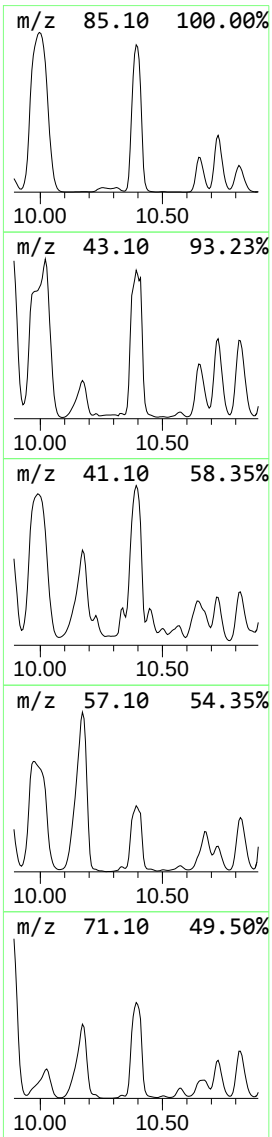
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Octane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.392	425.36 ug/l	86239600	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	95
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3	Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	50
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47
5	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

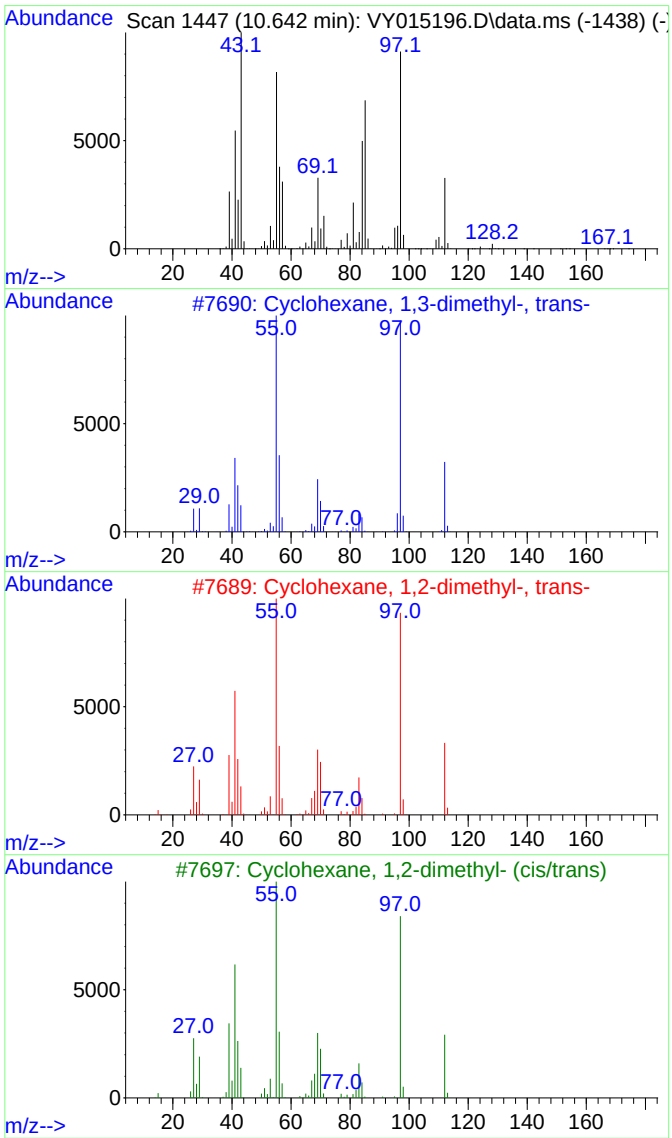
Quant Title :

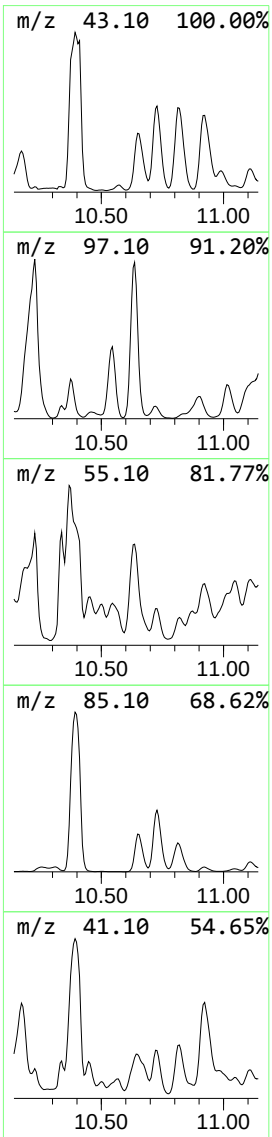
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.642	198.11 ug/l	40165700	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	64
2	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	53
3	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	53
4	Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	49
5	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	49





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

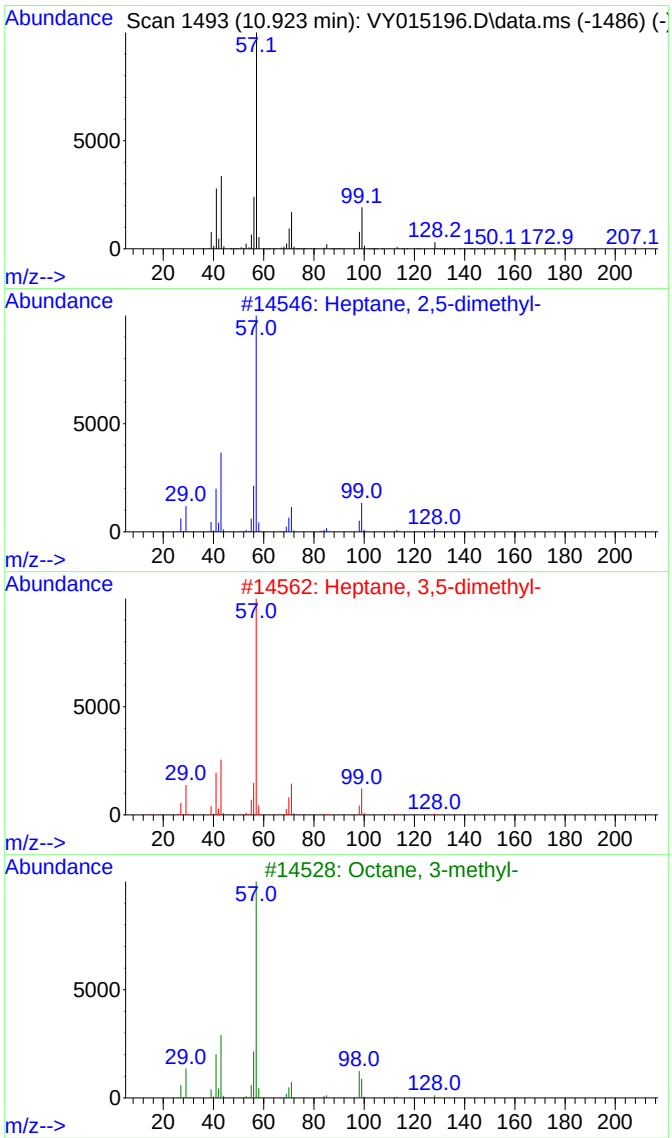
Quant Title :

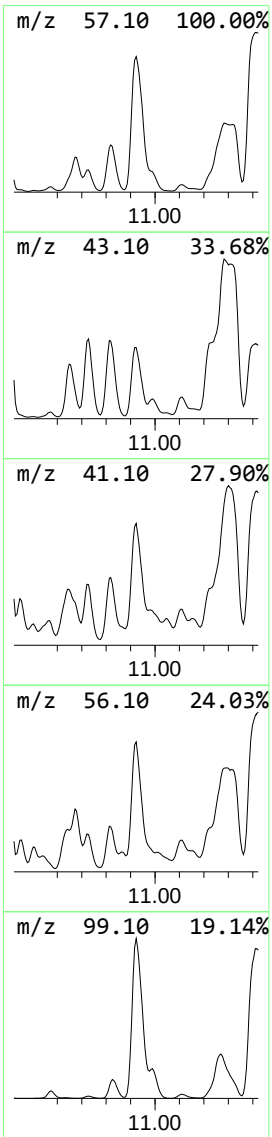
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.923	248.52 ug/l	50386400	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	91
2	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	87
3	Octane, 3-methyl-	128	C9H20	002216-33-3	59
4	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	53
5	Neopentyl ethyl ketone	128	C8H16O	005340-30-7	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

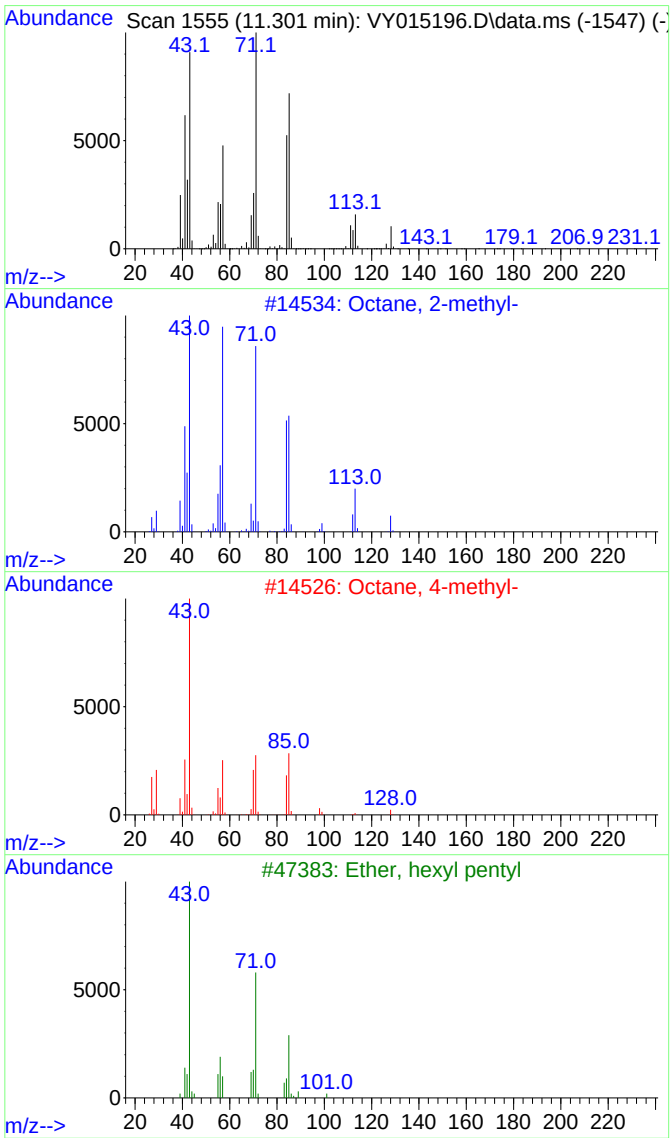
Quant Title :

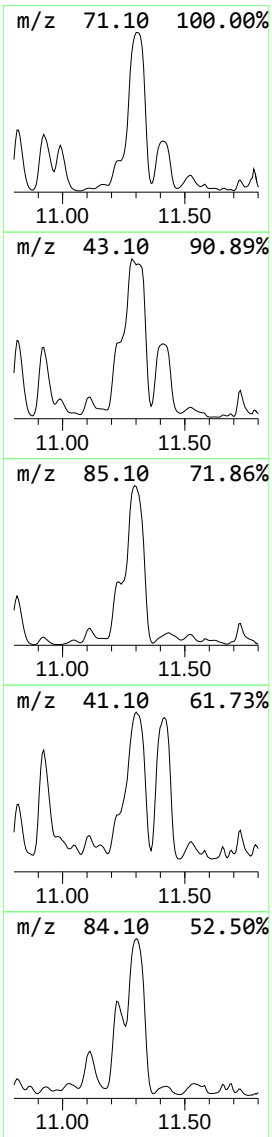
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.301	599.78 ug/l	121604000	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	81
2	Octane, 4-methyl-	128	C9H20	002216-34-4	72
3	Ether, hexyl pentyl	172	C11H24O	032357-83-8	50
4	Sulfurous acid, hexyl 2-pentyl e...	236	C11H24O3S	1000309-15-6	47
5	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

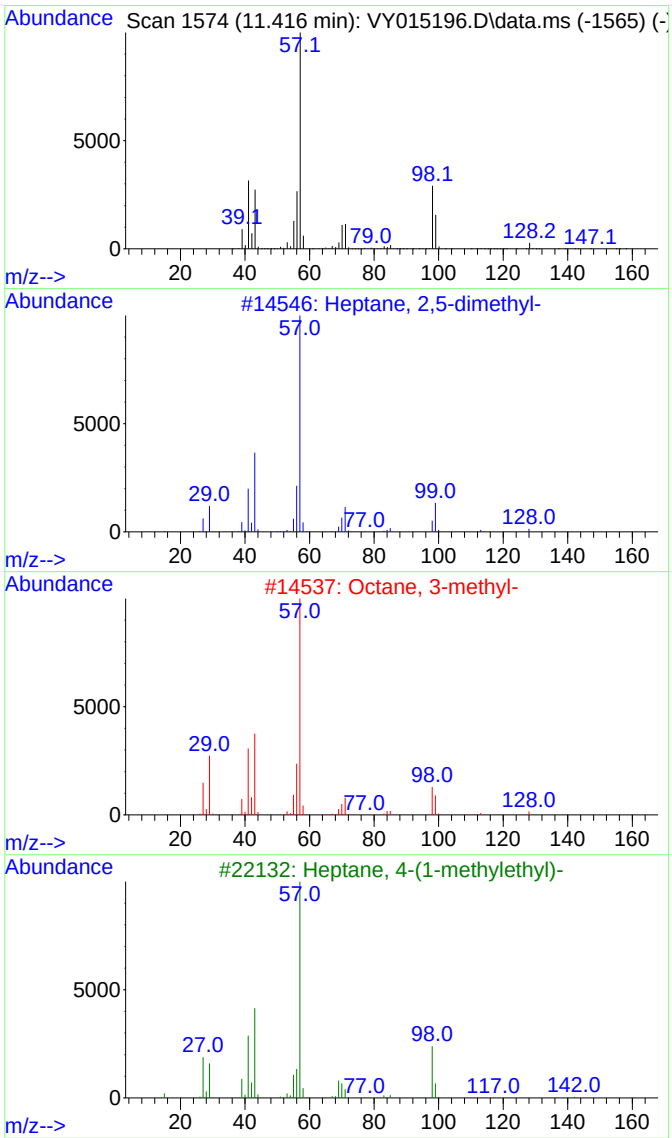
Quant Title :

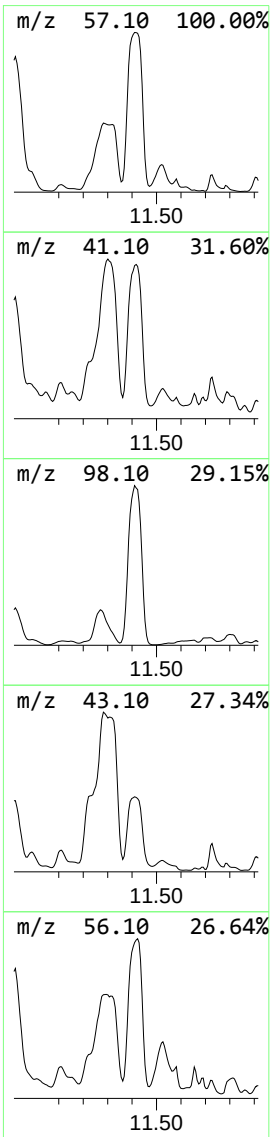
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.416	447.58 ug/l	90745400	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	74
2	Octane, 3-methyl-	128	C9H20	002216-33-3	72
3	Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	59
4	Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	53
5	Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	50





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Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

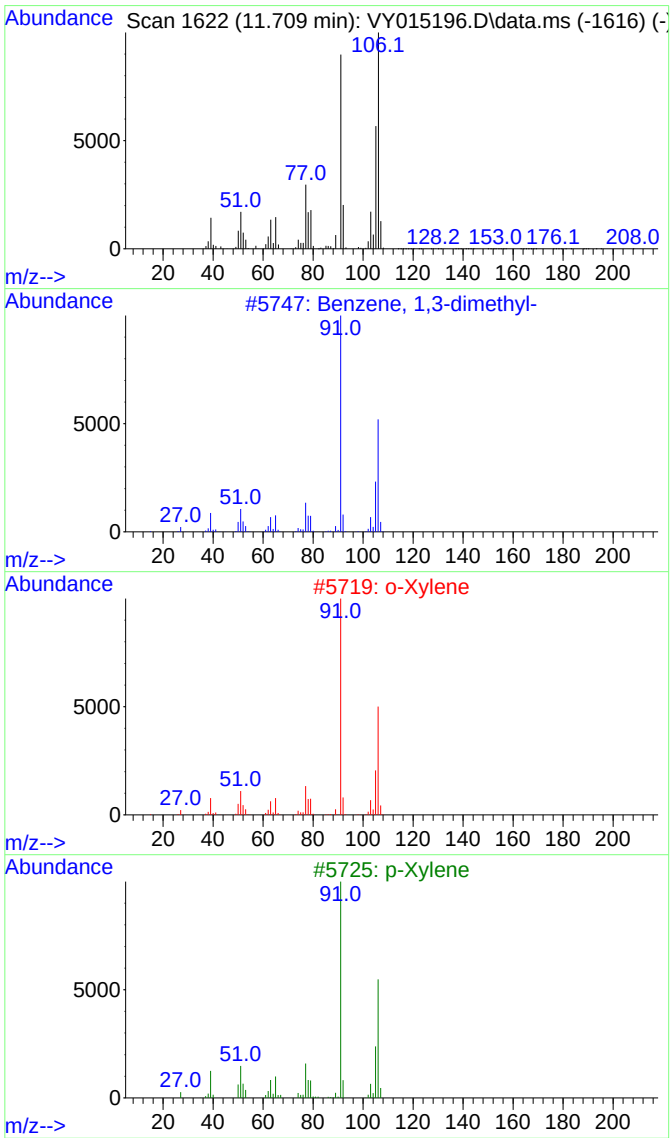
Quant Title :

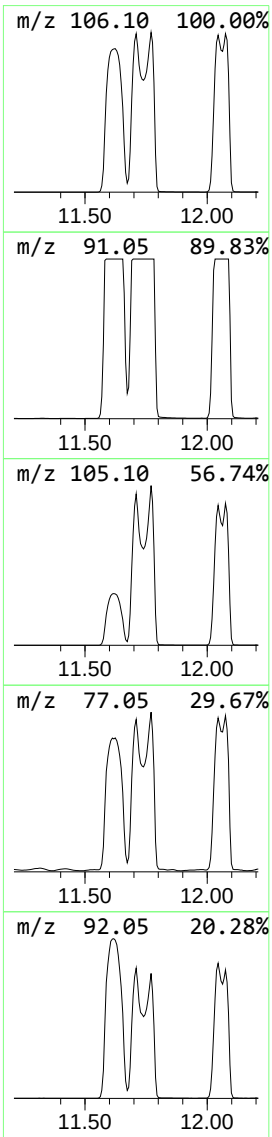
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.709	636.30 ug/l	129008000	Chlorobenzene-d5	11.502

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	93
2	o-Xylene	106	C8H10	000095-47-6	93
3	p-Xylene	106	C8H10	000106-42-3	90
4	Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	80
5	1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY081823\
Data File : VY015196.D
Acq On : 18 Aug 2023 15:24
Operator : SY/MD
Sample : 04086-07
Misc : 5.04g/5.2mL/MSVOA_Y/SOIL
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
B-PP09C

Quant Title :

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, 2,2-dim...	8.362	270.3	ug/l	209123000	2	8.697	38681400	50.0
unknown9.862	9.862	169.8	ug/l	131356000	2	8.697	38681400	50.0
Heptane, 3-methyl-	9.996	176.7	ug/l	136659000	2	8.697	38681400	50.0
Octane, 2,2,6-t...	10.173	256.1	ug/l	51925800	3	11.502	10137400	50.0
Octane	10.392	425.4	ug/l	86239600	3	11.502	10137400	50.0
Cyclohexane, 1,...	10.642	198.1	ug/l	40165700	3	11.502	10137400	50.0
Heptane, 2,5-di...	10.923	248.5	ug/l	50386400	3	11.502	10137400	50.0
Octane, 2-methyl-	11.301	599.8	ug/l	121604000	3	11.502	10137400	50.0
Octane, 3-methyl-	11.416	447.6	ug/l	90745400	3	11.502	10137400	50.0
Benzene, 1,3-di...	11.709	636.3	ug/l	129008000	3	11.502	10137400	50.0