LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

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_X\Method\82X121323W.M
Title : SW846 8260

Signal : TIC: VX039536.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.642	68	91	92	rBV7	10809	53518	5.99%	0.246%
2	5.391	692	706	714	rBV	61998	196345	21.98%	0.901%
3	5.489	715	722	725	rVV2	38219	104348	11.68%	0.479%
4	5.544	725	731	761	rVB2	106859	403288	45.14%	1.850%
5	5.812	761	775	787	rBV2	45392	135550	15.17%	0.622%
6	5.952	787	798	820	rBV	75032	222193	24.87%	1.019%
7	6.233	835	844	855	rVV2	104441	327818	36.70%	1.504%
8	6.598	893	904	914	rBV2	30279	75844	8.49%	0.348%
9	6.757	921	930	940	rBV	189719	490757	54.94%	2.251%
10	7.367	1019	1030	1040	rVB4	45198	126112	14.12%	0.579%
11	7.488	1042	1050	1055	rBV2	42463	90087	10.08%	0.413%
12	7.549	1055	1060	1072	rVB	52513	114200	12.78%	0.524%
13	7.976	1120	1130	1142	rVB	123920	272550	30.51%	1.250%
14	8.098	1142	1150	1159	rBV2	110661	263089	29.45%	1.207%
15	8.183	1159	1164	1173	rVB2	68204	129061	14.45%	0.592%
16	8.269	1173	1178	1182	rBV	137794	236346	26.46%	1.084%
17	8.305	1182	1184	1190	rVB	73047	103546	11.59%	0.475%
18	8.433	1200	1205	1213	rBV	155927	279073	31.24%	1.280%
19	8.507	1213	1217	1226	rVB4	25112	57289	6.41%	0.263%
20	8.610	1226	1234	1236	rBV4	93502	193747	21.69%	0.889%
21	8.647	1236	1240	1249	rVB	398951	662348	74.14%	3.039%
22	8.769	1249	1260	1264	rBV5	39562	93673	10.49%	0.430%
23	8.909	1277	1283	1290	rBV3	70979	141961	15.89%	0.651%
24	9.098	1306	1314	1318	rBV2	41386	88152	9.87%	0.404%
25	9.159	1320	1324	1327	rVV2	45589	73768	8.26%	0.338%
26	9.195	1327	1330	1340	rVB2	30055	53418	5.98%	0.245%
27	9.287	1340	1345	1349	rBV	36304	49876	5.58%	0.229%
28	9.378	1355	1360	1366	rBV2	55880	98292	11.00%	0.451%
29	9.494	1374	1379	1386	rBV3	108074	210410	23.55%	0.965%
30	9.653	1402	1405	1416	rVB8	30975	88043	9.86%	0.404%
31	9.762	1416	1423	1428	rBV5	47213	104047	11.65%	0.477%
32	9.817	1428	1432	1436	rBV2	41104	81202	9.09%	0.373%
33	9.866	1437	1440	1442	rVV3	41226	63433	7.10%	0.291%
34	9.903	1442	1446	1451	rVB	317651	433888	48.57%	1.990%
35	10.006	1458	1463	1467	rBV	253241	325062	36.39%	1.491%
36	10.055	1467	1471	1476	rVV	408633	575755	64.45%	2.641%

LSC Area Percent Report

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Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

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_X\Method\82X121323W.M
Title : SW846 8260

37	10.147	1481	1486	1489	rVV4	23614	52973	5.93%	0.243%
38	10.195	1490	1494	1498	rVB	51587	79125	8.86%	0.363%
39	10.262	1498	1505	1510	rBV3	53525	139689	15.64%	0.641%
40	10.360	1516	1521	1532	rVB8	77342	231967	25.97%	1.064%
41	10.585	1553	1558	1561	rBV2	90323	115394	12.92%	0.529%
42	10.671	1567	1572	1577	rBV2	78353	129828	14.53%	0.596%
43	10.781	1582	1590	1593	rBV	104064	183732	20.57%	0.843%
44	10.854	1593	1602	1607	rBV3	78407	141129	15.80%	0.647%
45	10.963	1613	1620	1627	rBV	96638	175113	19.60%	0.803%
46	11.024	1627	1630	1634	rBV2	63346	95727	10.72%	0.439%
47	11.085	1634	1640	1642	rVV	458116	691041	77.35%	3.170%
48	11.110	1642	1644	1651	rVB2	262930	359642	40.26%	1.650%
49	11.189	1651	1657	1661	rBV	184809	232035	25.97%	1.064%
50	11.305	1672	1676	1683	rVB	327549	407906	45.66%	1.871%
51	11.476	1700	1704	1711	rVB3	94948	162095	18.14%	0.744%
52	11.628	1723	1729	1734	rBV5	50557	99301	11.12%	0.456%
53	11.683	1734	1738	1740	rBV2	47755	57323	6.42%	0.263%
54	11.713	1740	1743	1746	rBV	52085	49287	5.52%	0.226%
55	11.750	1746	1749	1753	rBV2	47566	78506	8.79%	0.360%
56	11.835	1759	1763	1766	rBV	66589	99209	11.11%	0.455%
57	11.866	1766	1768	1770	rVV	55726	68556	7.67%	0.315%
58	11.890	1770	1772	1775	rVV	78869	86083	9.64%	0.395%
59	11.933	1775	1779	1783	rVV2	74097	125653	14.07%	0.576%
60	11.969	1783	1785	1788	rVV2	43578	50968	5.71%	0.234%
61	12.024	1788	1794	1799	rVV2	473859	735155	82.29%	3.373%
62	12.073	1799	1802	1804	rVV	112419	154489	17.29%	0.709%
63	12.103	1804	1807	1809	rVV	126248	176315	19.74%	0.809%
64	12.146	1812	1814	1816	rVV2	77924	95228	10.66%	0.437%
65	12.171	1816	1818	1825	rVB2	120645	145163	16.25%	0.666%
66	12.244	1825	1830	1836	rBV2	530586	718610	80.44%	3.297%
67	12.311	1836	1841	1849	rVB2	204882	465570	52.12%	2.136%
68	12.421	1853	1859	1863	rVV2	69298	128795	14.42%	0.591%
69	12.469	1863	1867	1873	rVB3	90431	141926	15.89%	0.651%
70	12.530	1873	1877	1883	rBV	53396	106289	11.90%	0.488%
71	12.616	1886	1891	1894	rVV	680677	886332	99.22%	4.066%
72	12.640	1894	1895	1900	rVV	99027	119634	13.39%	0.549%
73	12.701	1900	1905	1914	rVB2	264119	497277	55.66%	2.281%
74	12.920	1933	1941	1945	rBV	433314	655069	73.33%	3.005%
75	12.963	1945	1948	1953	rVB2	167338	252024	28.21%	1.156%
76	13.024	1953	1958	1967	rBV3	128129	284281	31.82%	1.304%

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Instrument :
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ClientSampleId :
SB32-2-20231222-16.0-17.0

Integration Parameters: RTEINT.P
Integrator: RTE
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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Title : SW846 8260

77	13.170	1978	1982	1985	rVV	308264	378702	42.39%	1.737%
78	13.213	1985	1989	1992	rVB3	47451	57786	6.47%	0.265%
79	13.262	1992	1997	1998	rBV2	100237	124186	13.90%	0.570%
80	13.292	1998	2002	2008	rVV2	621035	893341	100.00%	4.098%
81	13.372	2011	2015	2020	rVB2	132286	182337	20.41%	0.836%
82	13.420	2020	2023	2027	rBV2	52598	72071	8.07%	0.331%
83	13.506	2032	2037	2040	rBV2	96694	148881	16.67%	0.683%
84	13.542	2040	2043	2046	rVV	203475	275591	30.85%	1.264%
85	13.579	2046	2049	2055	rVV3	252270	435880	48.79%	2.000%
86	13.664	2056	2063	2070	rVV2	170017	461382	51.65%	2.117%
87	13.731	2070	2074	2077	rVV3	43699	89831	10.06%	0.412%
88	13.774	2077	2081	2087	rVV	169317	275887	30.88%	1.266%
89	13.829	2087	2090	2099	rVB5	60262	141921	15.89%	0.651%
90	13.926	2099	2106	2110	rBV3	114099	195301	21.86%	0.896%
91	13.969	2110	2113	2117	rVV	75332	103167	11.55%	0.473%
92	14.073	2126	2130	2135	rBV	166473	201460	22.55%	0.924%
93	14.213	2148	2153	2163	rBV	151823	252281	28.24%	1.157%
94	14.317	2167	2170	2176	rBV	75847	108485	12.14%	0.498%
95	14.371	2176	2179	2183	rVB	97881	113826	12.74%	0.522%
96	14.481	2192	2197	2202	rBV4	27781	63258	7.08%	0.290%
97	14.634	2214	2222	2227	rBV	304589	408713	45.75%	1.875%
98	14.780	2240	2246	2250	rBV	162917	211222	23.64%	0.969%
99	15.475	2355	2360	2375	rVV	31873	56057	6.27%	0.257%
100	15.615	2377	2383	2386	rVV2	33689	53891	6.03%	0.247%

Sum of corrected areas: 21797985

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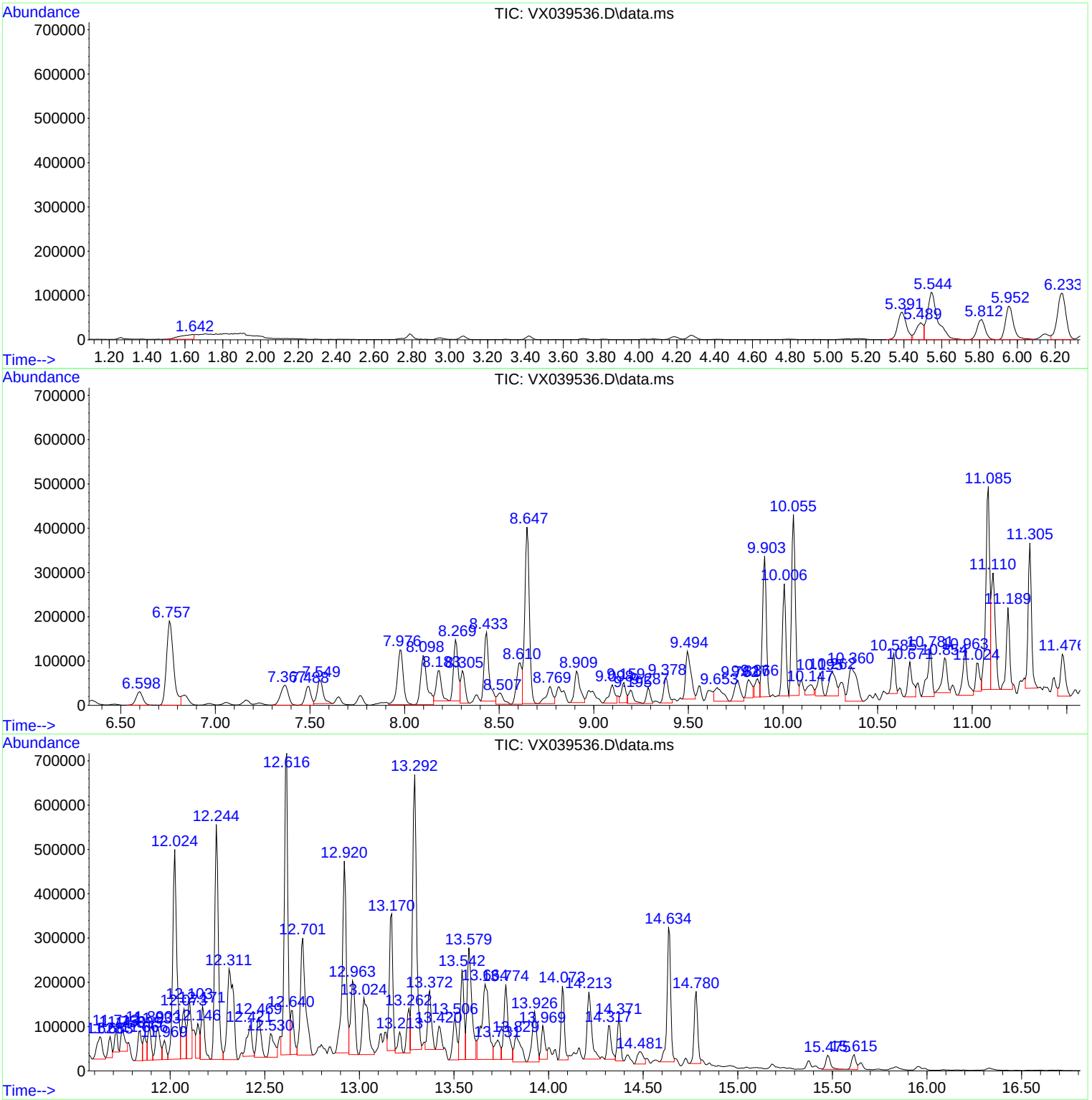
Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
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Instrument :
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ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

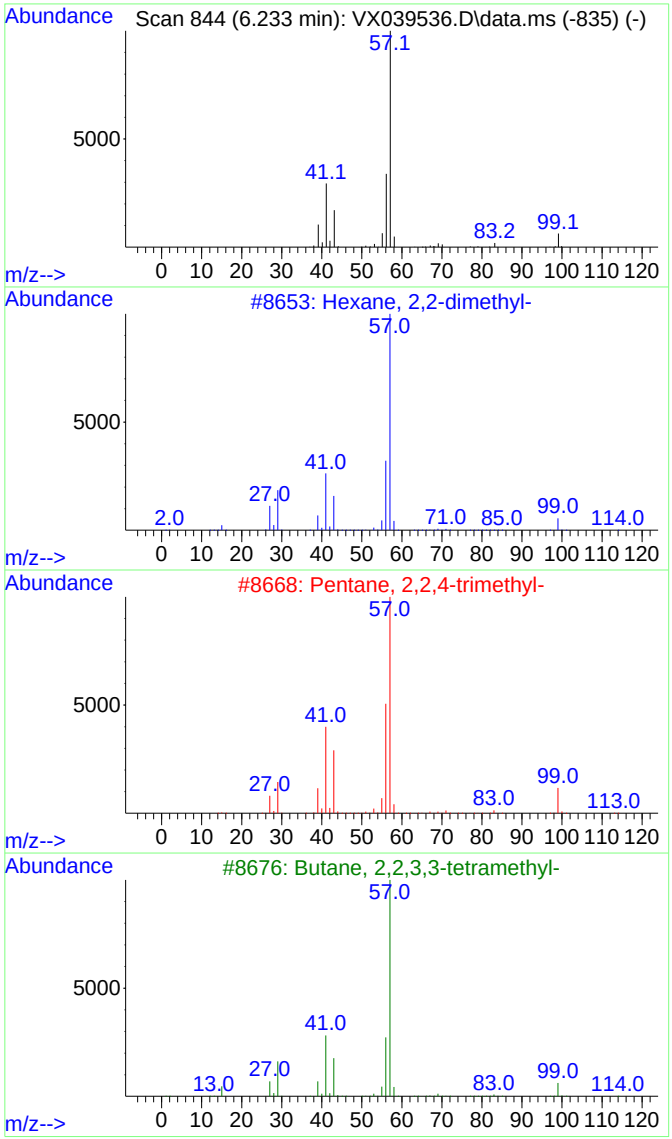
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

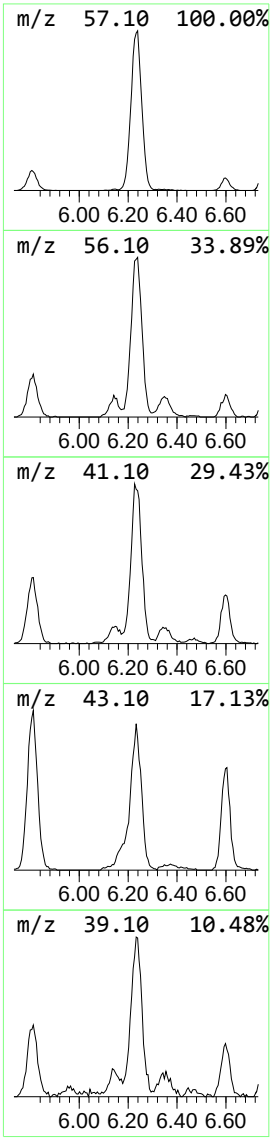
****

Peak Number 1 Hexane, 2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.233	33.40 ug/l	327818	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
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Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

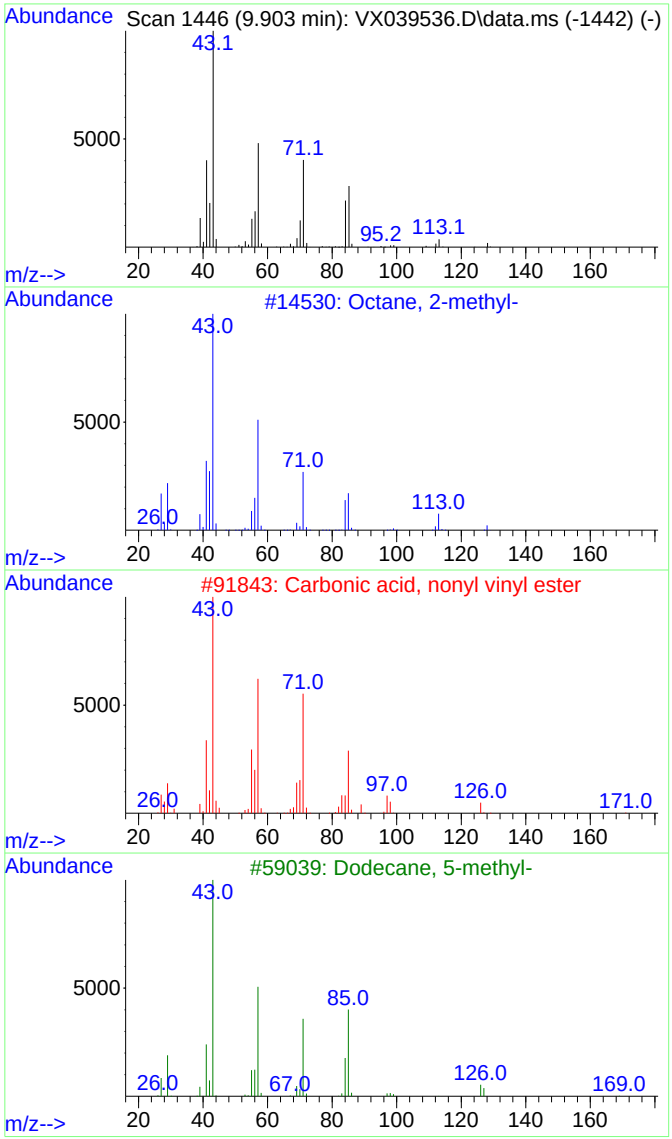
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

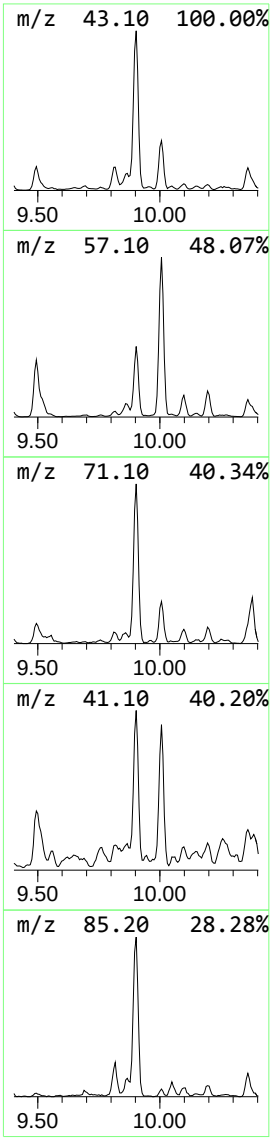
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.903	37.68 ug/l	433888	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	87
2	Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	64
3	Dodecane, 5-methyl-	184	C13H28	017453-93-9	59
4	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
5	Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	53





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Instrument :
MSVOA_X
ClientSampleId :
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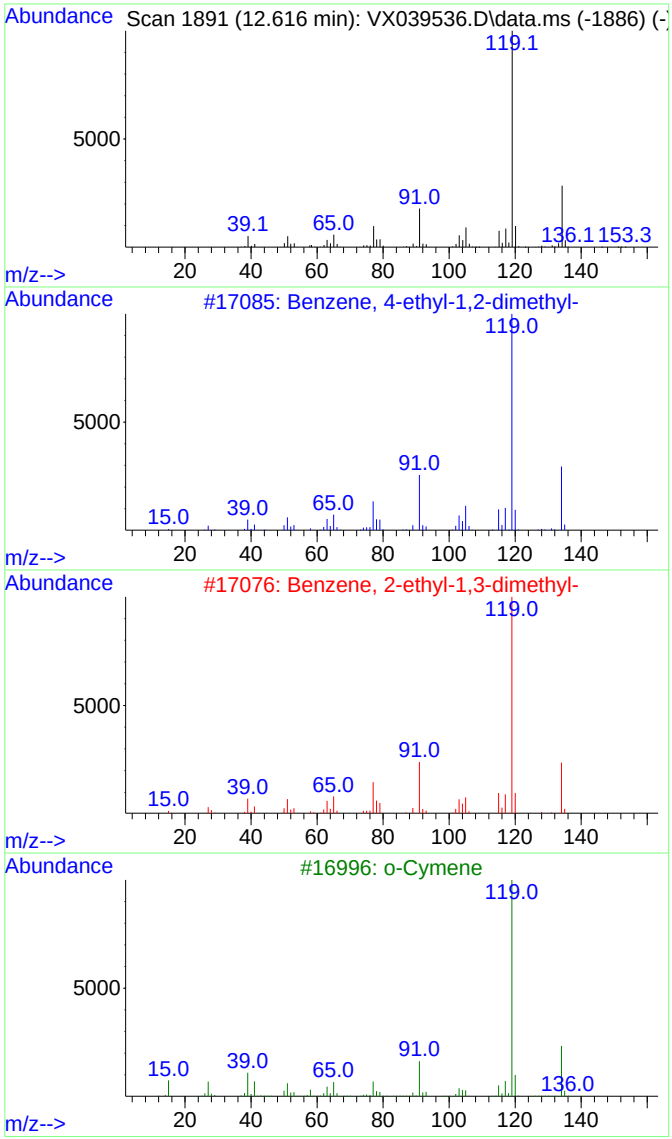
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

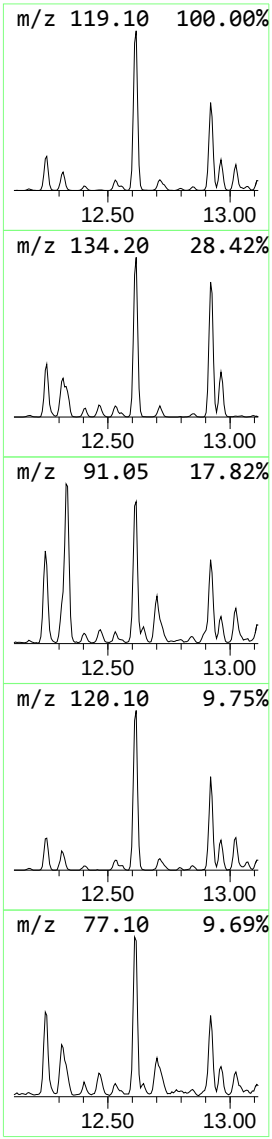
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	60.28 ug/l	886332	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





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Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

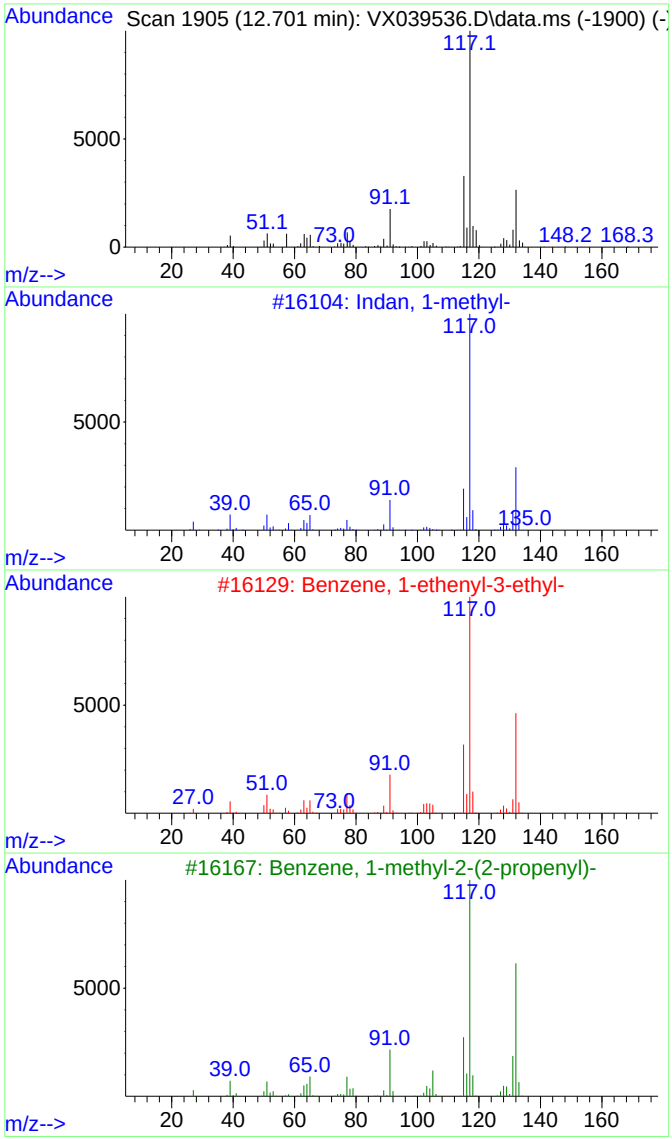
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

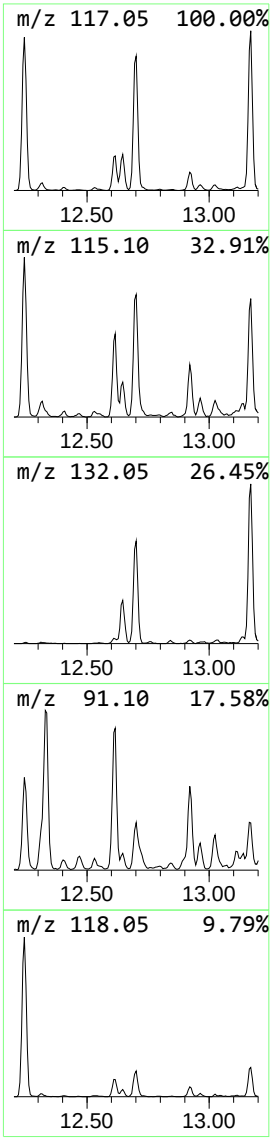
****

Peak Number 5 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	33.82 ug/l	497277	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

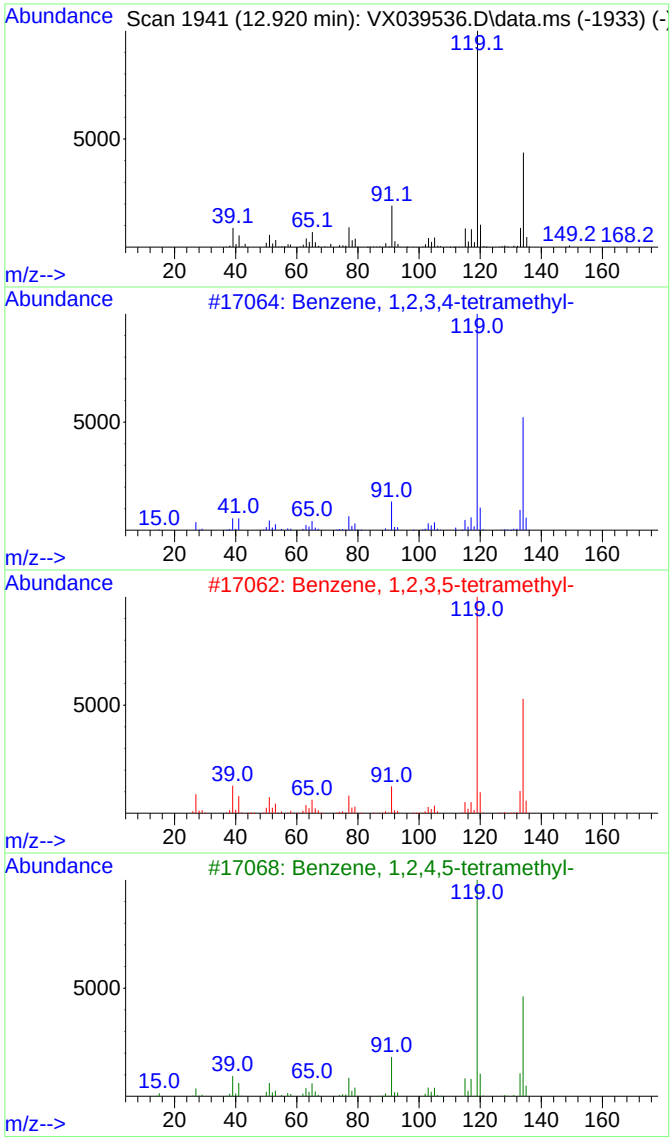
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

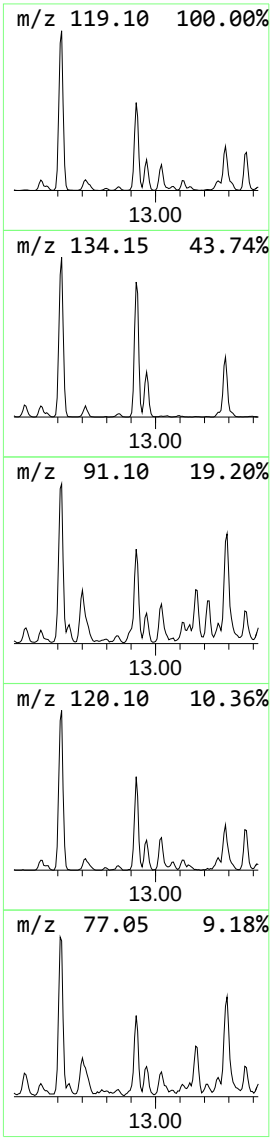
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	44.55 ug/l	655069	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94





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Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

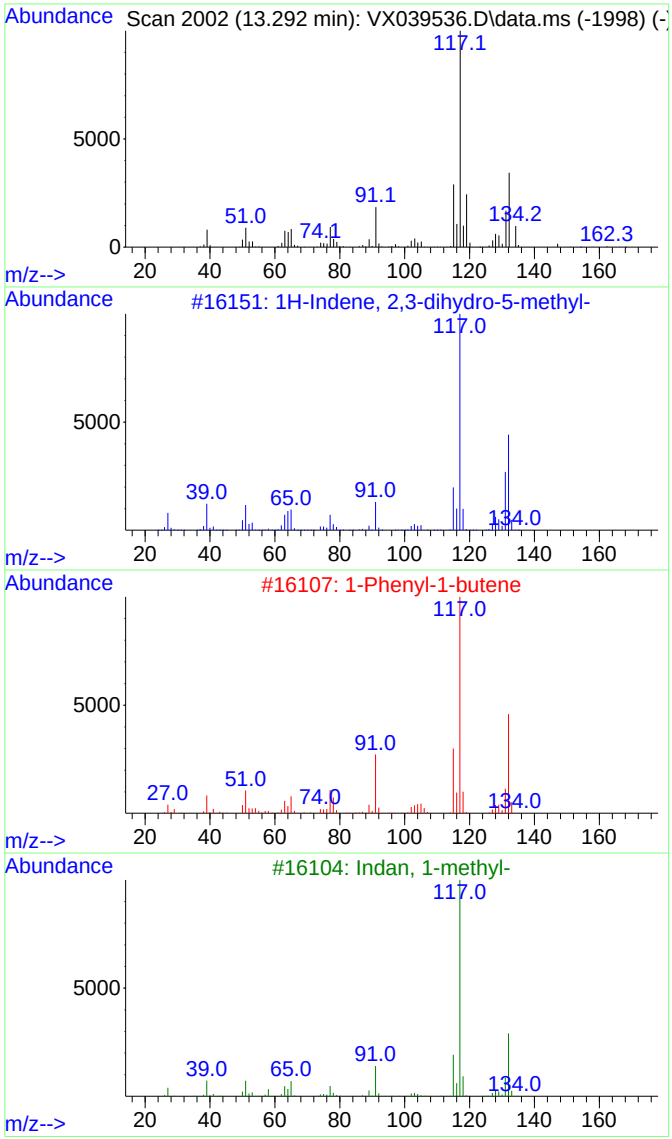
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

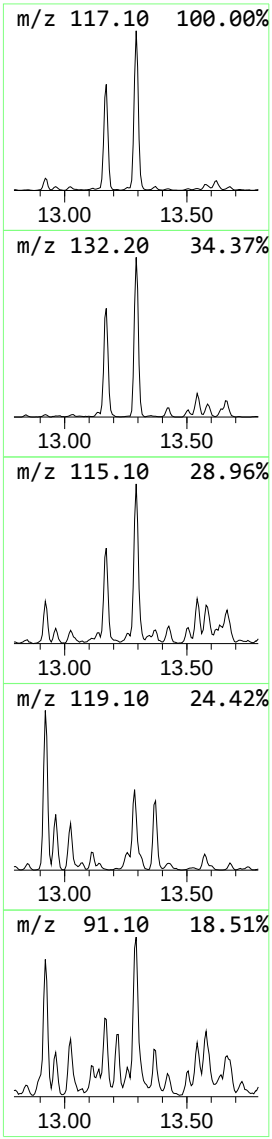
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	60.76 ug/l	893341	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

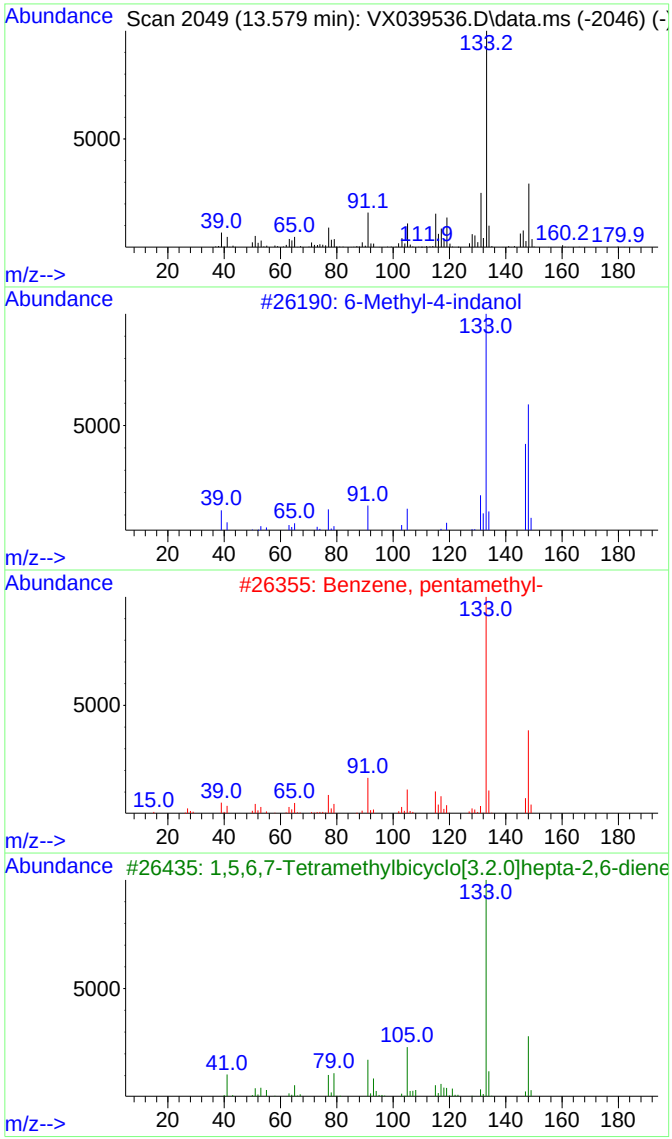
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

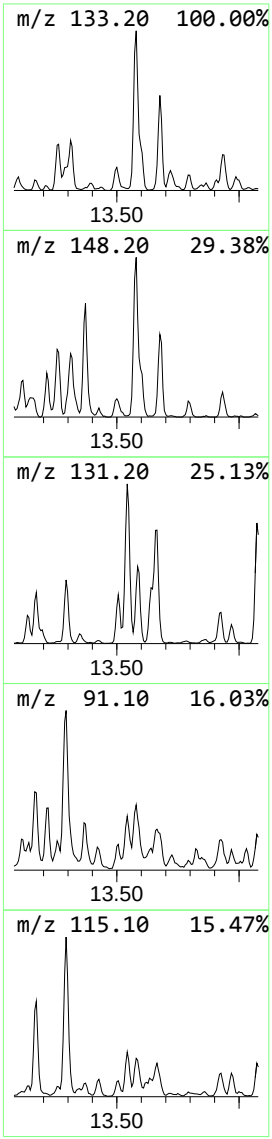
****

Peak Number 8 6-Methyl-4-indanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	29.65 ug/l	435880	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	80
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	76
3	1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	68
4	3,4-Dimethylcumene	148	C11H16	1000370-34-1	64
5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

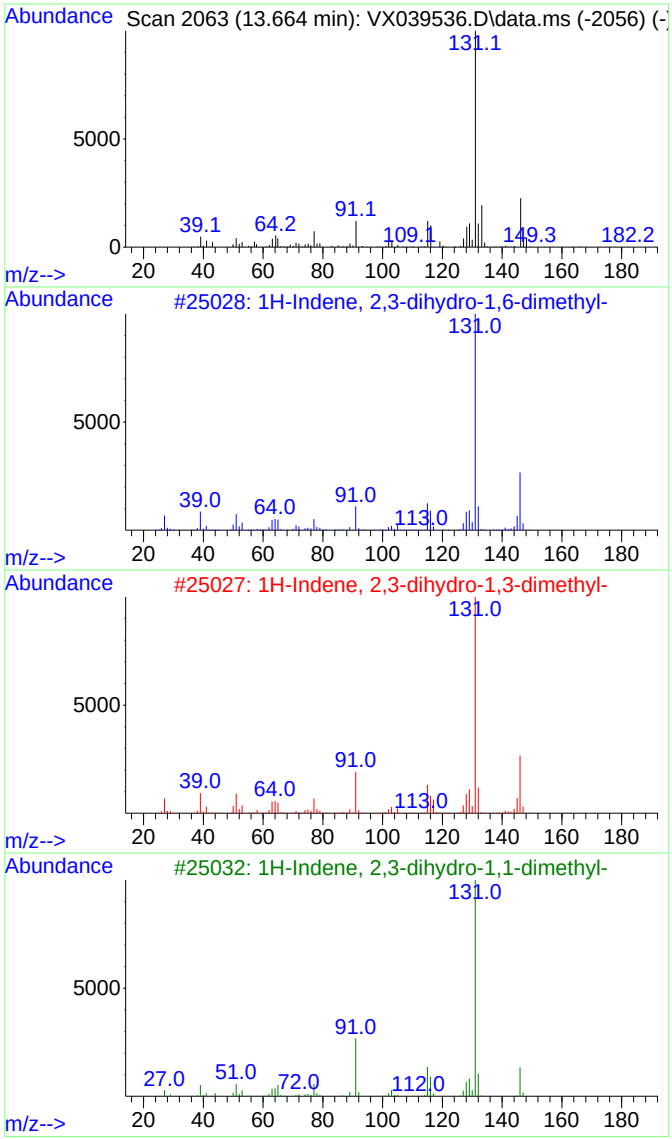
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

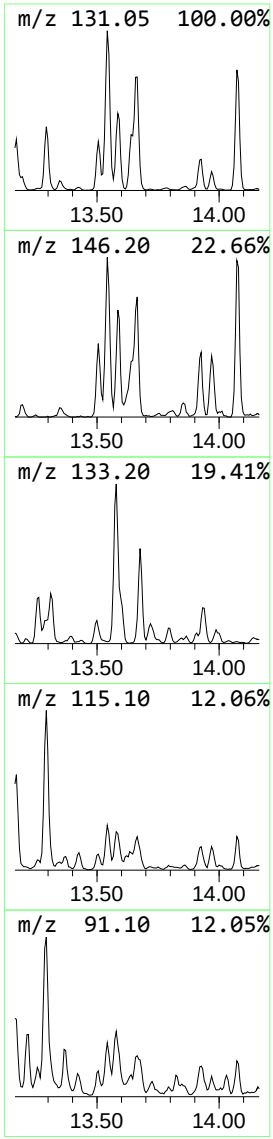
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	31.38 ug/l	461382	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
3	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	87
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	87





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Operator : JC/MD
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Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

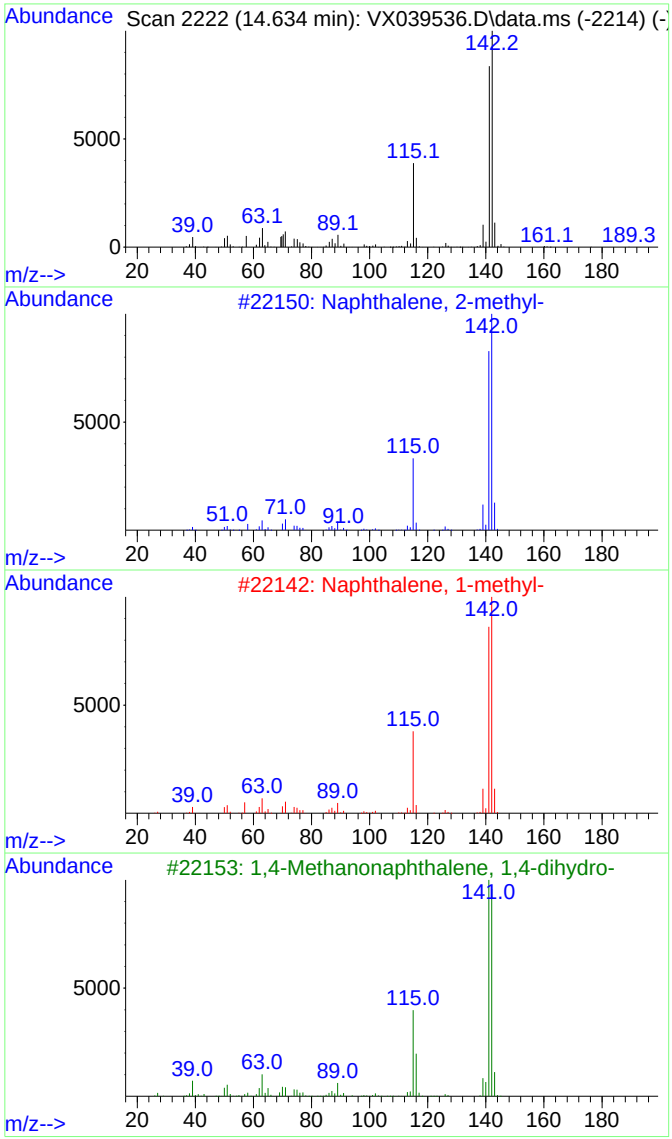
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
Quant Title : SW846 8260

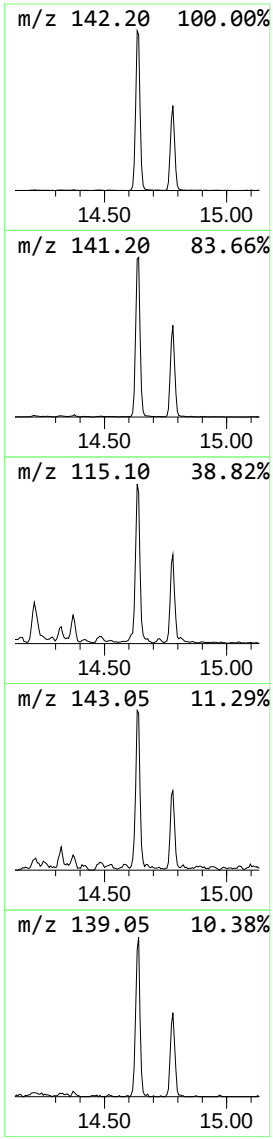
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	27.80 ug/l	408713	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
Data File : VX039536.D
Acq On : 29 Dec 2023 11:42
Operator : JC/MD
Sample : 06017-04
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB32-2-20231222-16.0-17.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.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, 2,2-dim...	6.233	33.4	ug/l	327818	2	6.757	490757	50.0
Octane, 2-methyl-	9.903	37.7	ug/l	433888	3	10.055	575755	50.0
Benzene, 4-ethy...	12.616	60.3	ug/l	886332	4	12.024	735155	50.0
Indan, 1-methyl-	12.701	33.8	ug/l	497277	4	12.024	735155	50.0
Benzene, 1,2,3,...	12.920	44.5	ug/l	655069	4	12.024	735155	50.0
1H-Indene, 2,3-...	13.292	60.8	ug/l	893341	4	12.024	735155	50.0
6-Methyl-4-indanol	13.579	29.6	ug/l	435880	4	12.024	735155	50.0
1H-Indene, 2,3-...	13.664	31.4	ug/l	461382	4	12.024	735155	50.0
1,4-Methanonaph...	14.634	27.8	ug/l	408713	4	12.024	735155	50.0