LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

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\82N030524W.M
Title : SW846 8260

Signal : TIC: VN081360.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.801	136	144	145	rBV2	35050	57744	1.06%	0.075%
2	8.165	1044	1056	1060	rBV2	260806	614451	11.24%	0.798%
3	8.218	1060	1065	1077	rVB	509901	1135063	20.76%	1.473%
4	8.577	1116	1126	1137	rBV	307286	676517	12.37%	0.878%
5	8.759	1142	1157	1166	rBV	69633	195400	3.57%	0.254%
6	9.100	1206	1215	1223	rBV	729879	1476468	27.00%	1.917%
7	9.594	1287	1299	1305	rBV3	90202	255053	4.66%	0.331%
8	9.724	1315	1321	1327	rBV	30871	70342	1.29%	0.091%
9	9.865	1335	1345	1353	rVB6	23815	64552	1.18%	0.084%
10	10.012	1361	1370	1379	rVB	356037	710561	12.99%	0.922%
11	10.147	1379	1393	1405	rBV	561266	1201023	21.96%	1.559%
12	10.324	1416	1423	1428	rBV4	45938	107269	1.96%	0.139%
13	10.371	1428	1431	1441	rVB5	32166	61348	1.12%	0.080%
14	10.494	1441	1452	1457	rBV	257221	480195	8.78%	0.623%
15	10.565	1457	1464	1477	rVB	1186896	2323513	42.49%	3.016%
16	10.747	1488	1495	1503	rVB5	64112	160257	2.93%	0.208%
17	10.912	1518	1523	1532	rVB5	81405	154438	2.82%	0.200%
18	11.012	1532	1540	1548	rBV4	184964	389616	7.12%	0.506%
19	11.082	1548	1552	1558	rVB7	33551	68443	1.25%	0.089%
20	11.235	1574	1578	1581	rBV3	46185	83286	1.52%	0.108%
21	11.353	1593	1598	1600	rBV5	39267	74607	1.36%	0.097%
22	11.382	1600	1603	1606	rVV5	67393	115711	2.12%	0.150%
23	11.418	1606	1609	1613	rVV2	105865	186634	3.41%	0.242%
24	11.471	1613	1618	1623	rVV2	173469	332760	6.08%	0.432%
25	11.524	1623	1627	1631	rVV5	66130	105075	1.92%	0.136%
26	11.571	1631	1635	1639	rVV5	70526	117653	2.15%	0.153%
27	11.612	1639	1642	1648	rVB3	72179	108830	1.99%	0.141%
28	11.676	1648	1653	1660	rBV4	118573	252873	4.62%	0.328%
29	11.865	1676	1685	1691	rBV	1001598	2000276	36.58%	2.597%
30	12.000	1705	1708	1712	rVB3	80674	109558	2.00%	0.142%
31	12.076	1712	1721	1724	rBV	181485	437655	8.00%	0.568%
32	12.112	1724	1727	1732	rVB2	221054	332001	6.07%	0.431%
33	12.218	1739	1745	1746	rBV6	43149	73178	1.34%	0.095%
34	12.306	1756	1760	1764	rVB5	52469	68819	1.26%	0.089%
35	12.400	1765	1776	1782	rBV3	359949	961799	17.59%	1.249%
36	12.471	1785	1788	1794	rVB3	121589	201168	3.68%	0.261%

LSC Area Percent Report

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Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

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\82N030524W.M
Title : SW846 8260

37	12.588	1794	1808	1816	rBV5	290924	1051923	19.24%	1.366%
38	12.829	1841	1849	1850	rBV	576623	962666	17.60%	1.250%
39	12.853	1850	1853	1860	rVB	835799	1336126	24.43%	1.734%
40	12.935	1860	1867	1872	rBV3	155335	306970	5.61%	0.398%
41	12.988	1872	1876	1879	rBV	120469	188511	3.45%	0.245%
42	13.029	1879	1883	1888	rVB3	152295	291104	5.32%	0.378%
43	13.100	1889	1895	1898	rBV	782509	1308679	23.93%	1.699%
44	13.135	1898	1901	1903	rVV	388910	569555	10.41%	0.739%
45	13.176	1903	1908	1915	rVB	1496072	2632429	48.14%	3.417%
46	13.341	1928	1936	1942	rBV	497120	936868	17.13%	1.216%
47	13.423	1943	1950	1954	rVB2	266308	497336	9.09%	0.646%
48	13.482	1954	1960	1966	rBV	3413951	5468712	100.00%	7.099%
49	13.535	1966	1969	1974	rVB2	139436	214769	3.93%	0.279%
50	13.647	1977	1988	1991	rBV3	690576	1526261	27.91%	1.981%
51	13.794	2008	2013	2015	rBV	1032865	1631681	29.84%	2.118%
52	13.823	2015	2018	2032	rVB2	1850740	3908148	71.46%	5.073%
53	13.953	2035	2040	2041	rBV	275050	351590	6.43%	0.456%
54	13.988	2041	2046	2047	rVV	1075473	1528048	27.94%	1.984%
55	14.017	2048	2051	2063	rVB	1916766	3553596	64.98%	4.613%
56	14.117	2063	2068	2074	rBV5	382565	777818	14.22%	1.010%
57	14.188	2075	2080	2085	rVB	536006	855552	15.64%	1.111%
58	14.247	2085	2090	2092	rBV	949168	1468400	26.85%	1.906%
59	14.276	2092	2095	2099	rVV	927969	1329192	24.31%	1.725%
60	14.335	2099	2105	2111	rVB	1737763	2792007	51.05%	3.624%
61	14.400	2111	2116	2118	rBV	169277	246316	4.50%	0.320%
62	14.447	2118	2124	2129	rVB2	986090	1633375	29.87%	2.120%
63	14.576	2140	2146	2151	rBV2	577063	980585	17.93%	1.273%
64	14.659	2151	2160	2163	rVV	995533	2150124	39.32%	2.791%
65	14.694	2163	2166	2170	rVV	1447606	2146185	39.24%	2.786%
66	14.735	2170	2173	2177	rVV	635533	976915	17.86%	1.268%
67	14.776	2177	2180	2188	rVB2	446144	867452	15.86%	1.126%
68	14.882	2189	2198	2201	rVV4	356129	743170	13.59%	0.965%
69	14.929	2201	2206	2210	rVV	966967	1799719	32.91%	2.336%
70	14.964	2210	2212	2218	rVB2	575590	859363	15.71%	1.116%
71	15.053	2218	2227	2233	rBV3	1448450	3955031	72.32%	5.134%
72	15.100	2233	2235	2247	rVV4	395348	813648	14.88%	1.056%
73	15.217	2250	2255	2262	rBV2	193573	414086	7.57%	0.538%
74	15.323	2262	2273	2286	rVV4	876189	2922950	53.45%	3.794%
75	15.441	2286	2293	2300	rVB5	684038	1685119	30.81%	2.188%
76	15.594	2315	2319	2322	rBV4	157127	273923	5.01%	0.356%

LSC Area Percent Report

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Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

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\82N030524W.M
Title : SW846 8260

77	15.641	2322	2327	2333	rVB	480171	896458	16.39%	1.164%
78	15.747	2341	2345	2350	rBV3	199580	345757	6.32%	0.449%
79	15.947	2372	2379	2388	rBV2	347021	842880	15.41%	1.094%
80	16.123	2402	2409	2413	rBV5	142155	370778	6.78%	0.481%
81	16.711	2504	2509	2515	rBV5	94815	212615	3.89%	0.276%
82	16.782	2515	2521	2522	rBV2	451503	645300	11.80%	0.838%

Sum of corrected areas: 77033826

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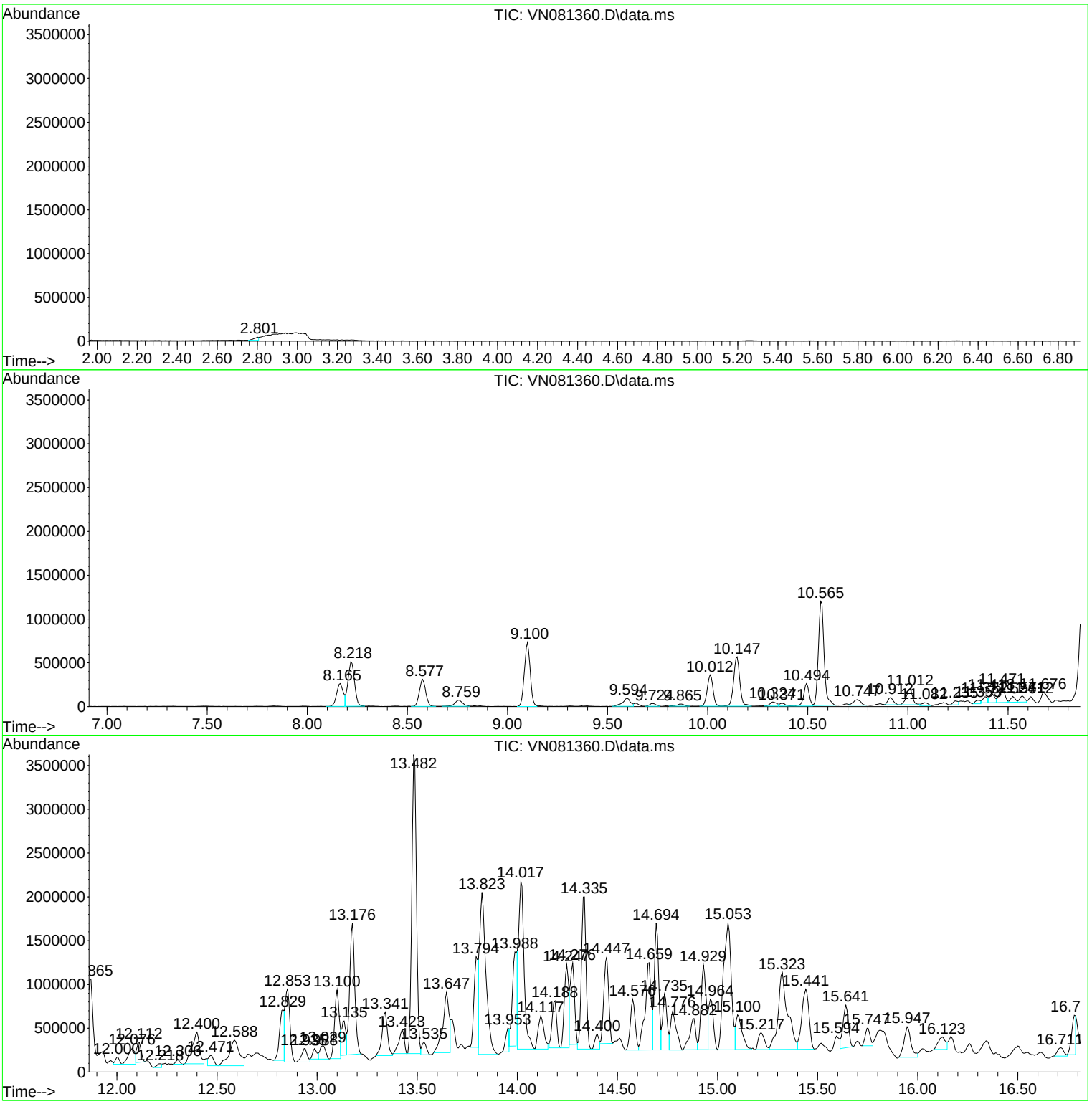
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

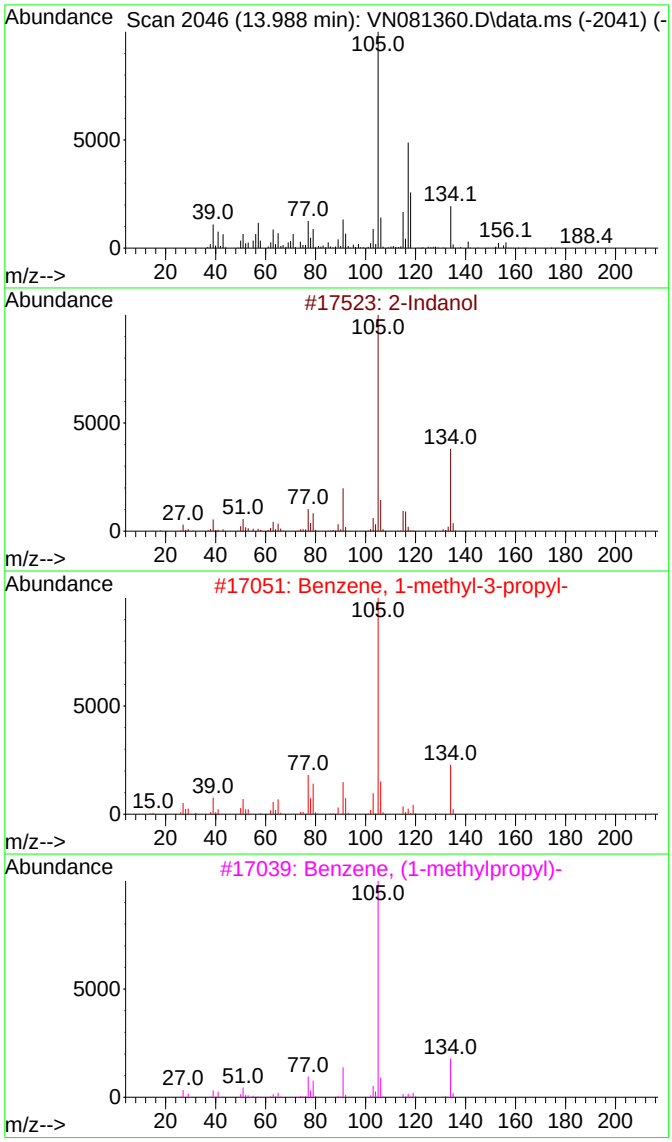
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

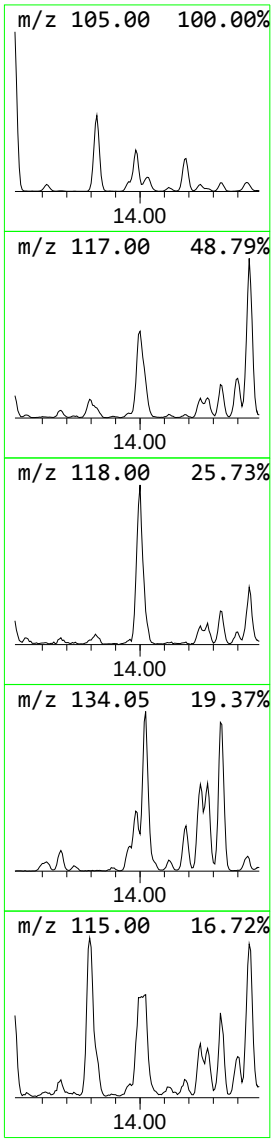
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 unknown13.988 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.988	46.82 ug/l	1528050	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Indanol	134	C9H10O	004254-29-9	38
2	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38
4	2-Tolyloxirane	134	C9H10O	002783-26-8	30
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

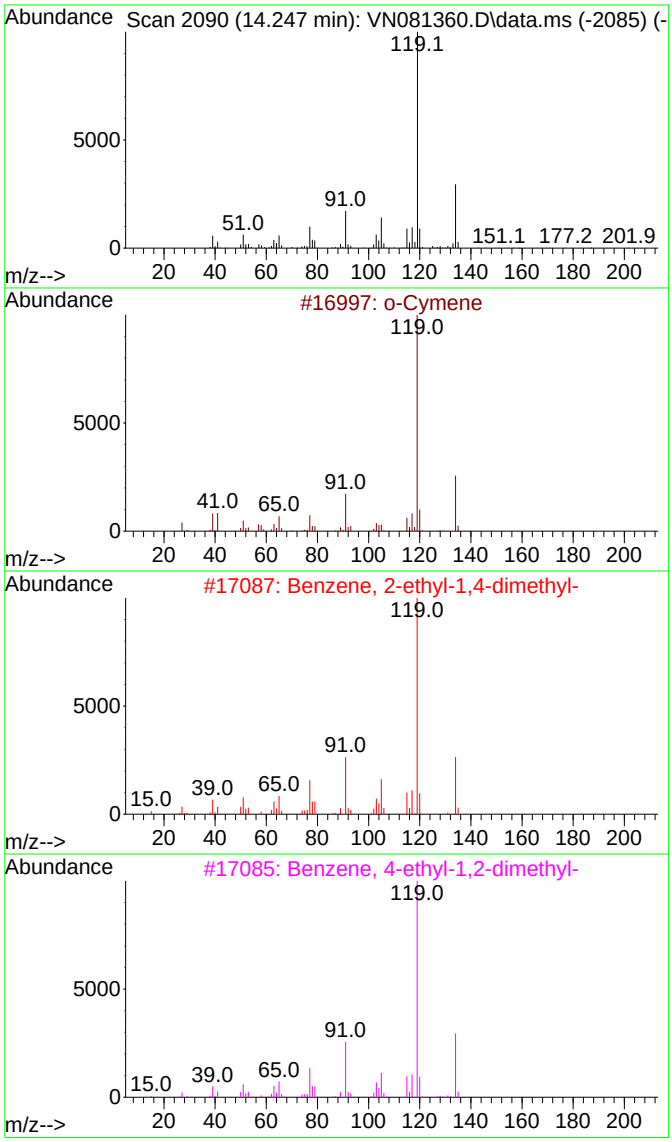
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

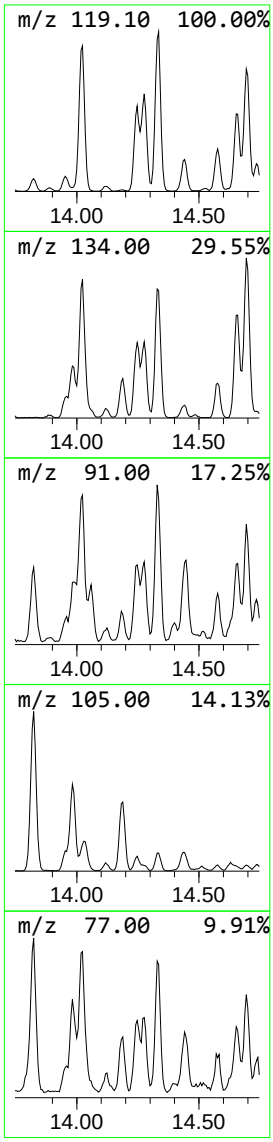
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.247	45.00 ug/l	1468400	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
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Operator : JC\MD
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Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

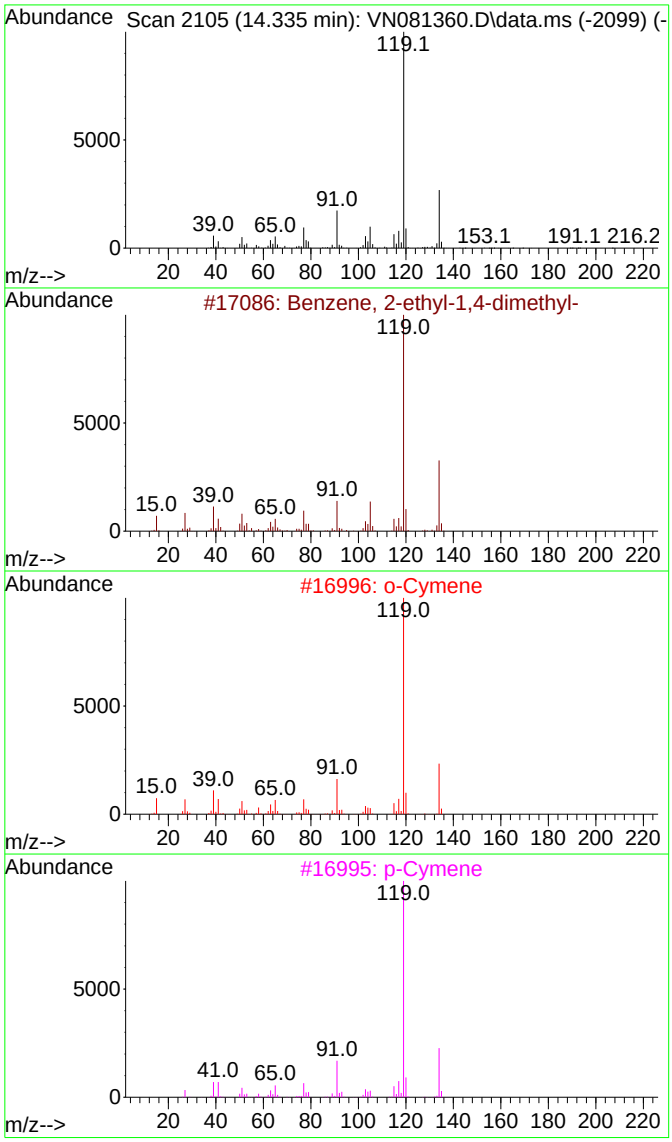
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

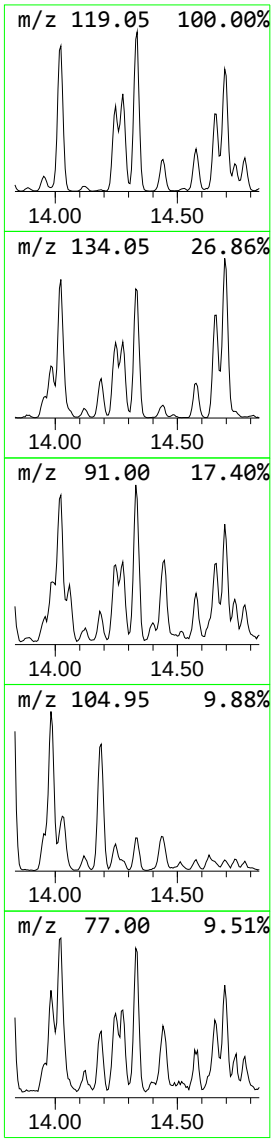
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.335	85.56 ug/l	2792010	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	p-Cymene	134	C10H14	000099-87-6	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93





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Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

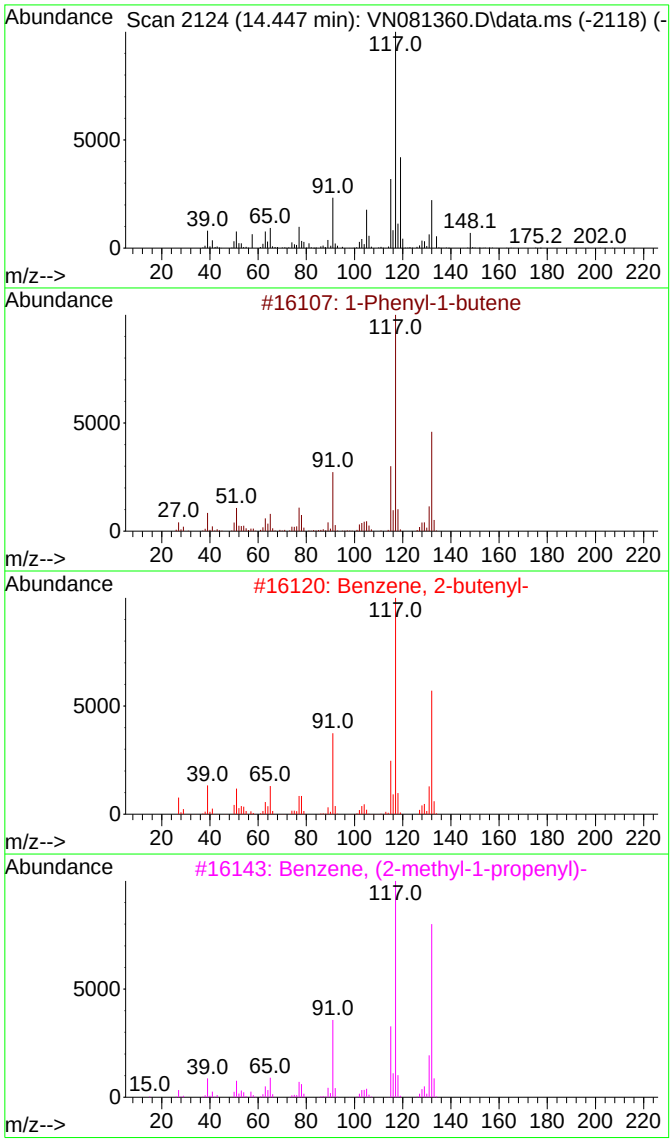
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

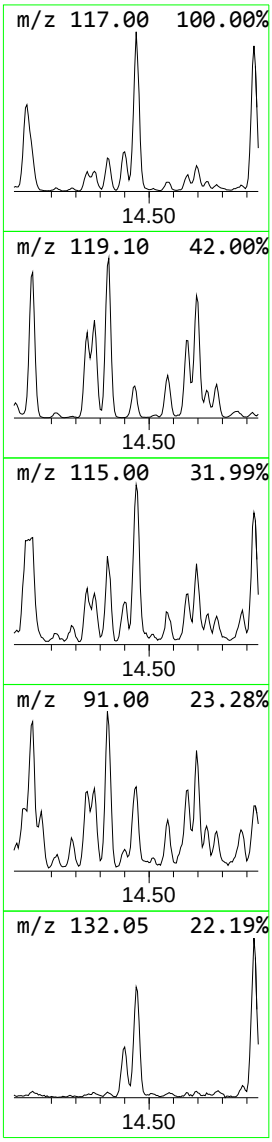
****

Peak Number 4 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	50.05 ug/l	1633380	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	62
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	62
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	62
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	60





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Data File : VN081360.D
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Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

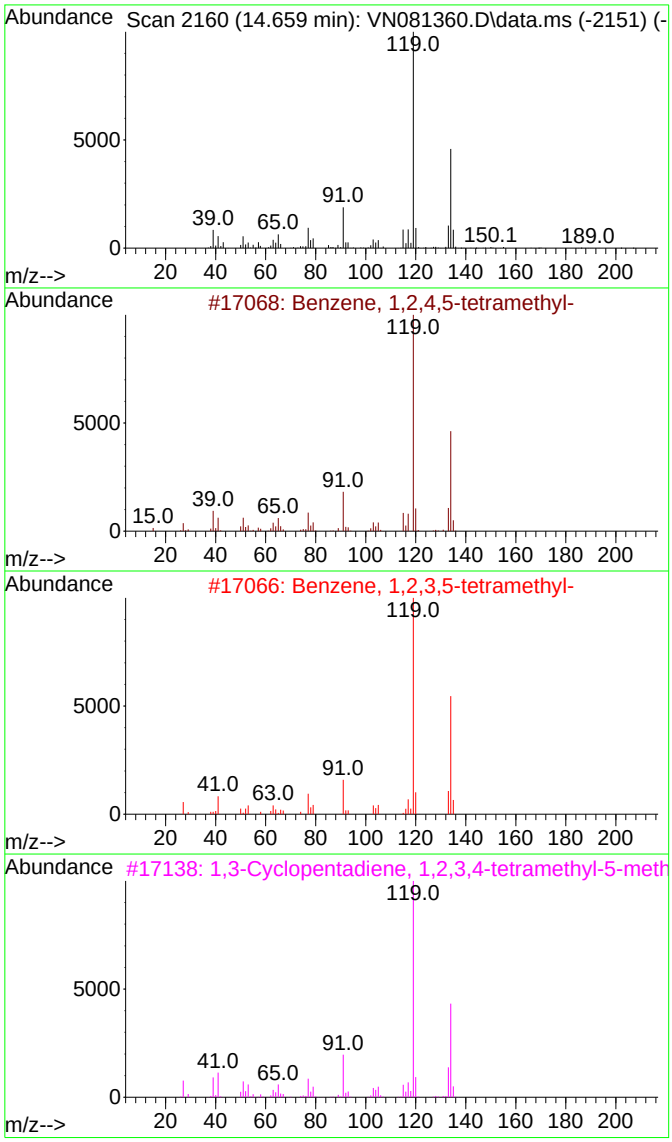
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

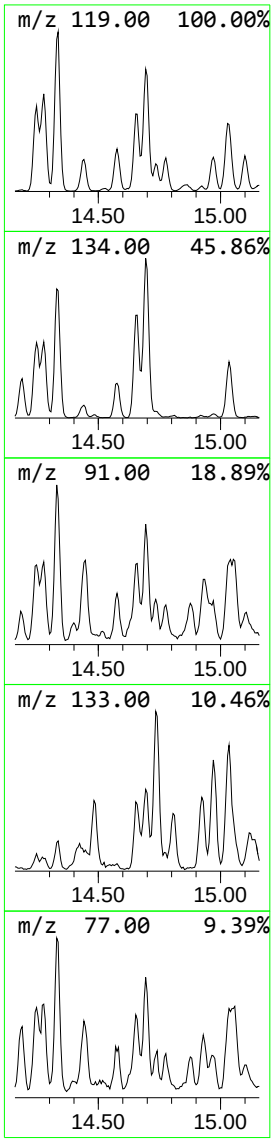
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.659	65.89 ug/l	2150120	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	95
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5	o-Cymene	134	C10H14	000527-84-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

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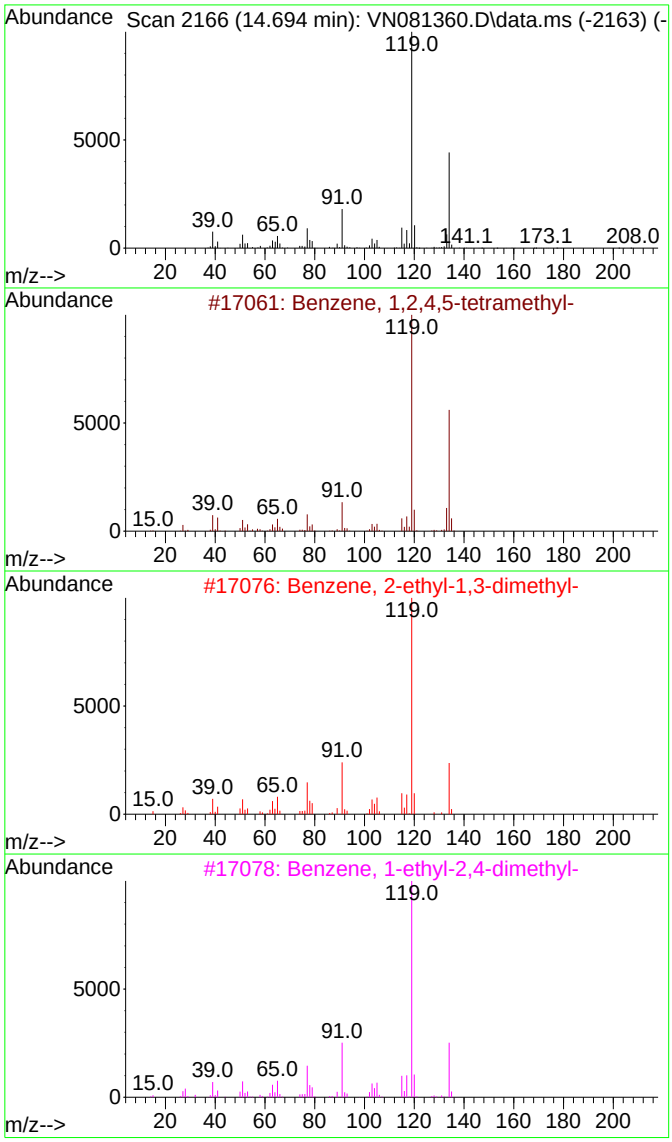
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

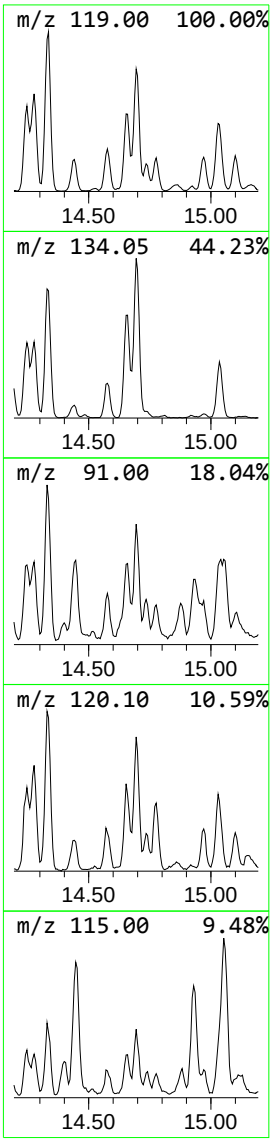
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.694	65.77 ug/l	2146190	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

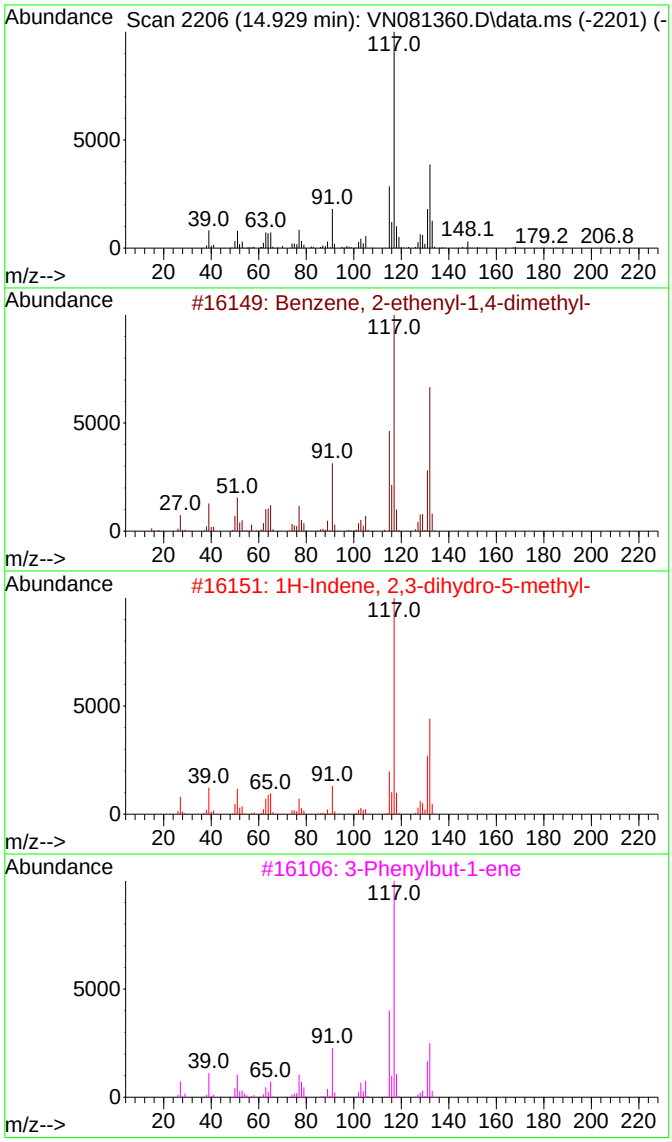
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

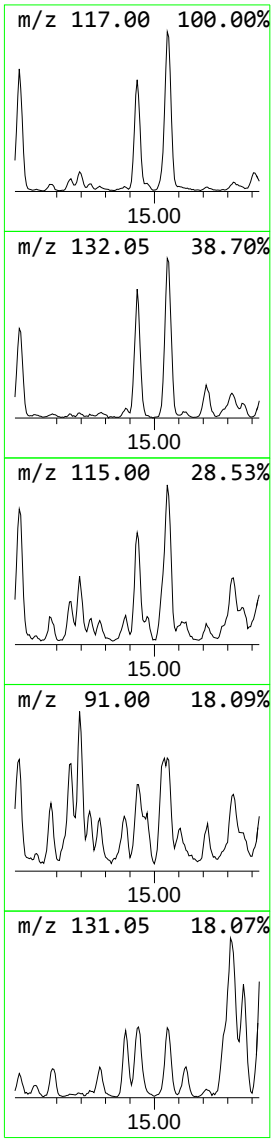
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.929	55.15 ug/l	1799720	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

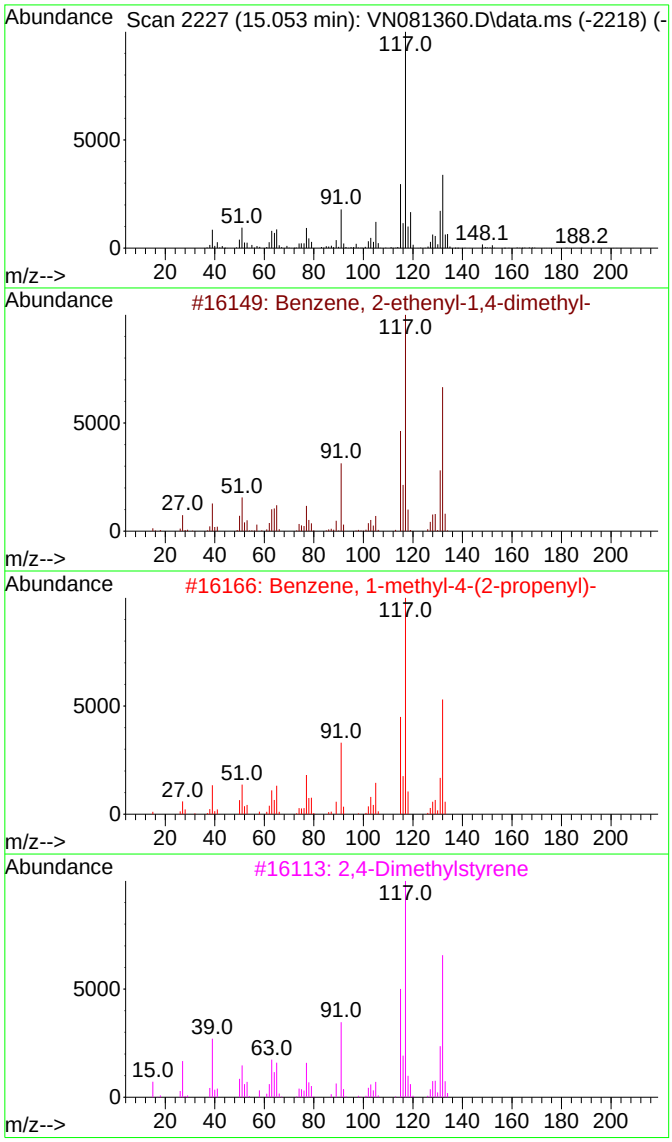
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

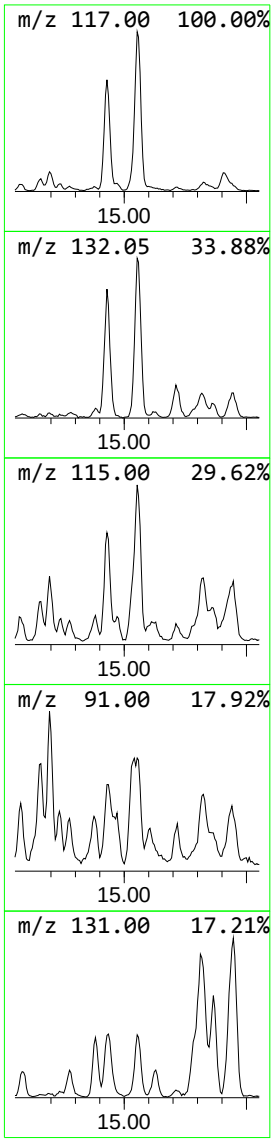
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-methyl-4-(2-prop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	121.19 ug/l	3955030	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	96
3	2,4-Dimethylstyrene	132	C10H12	002234-20-0	96
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

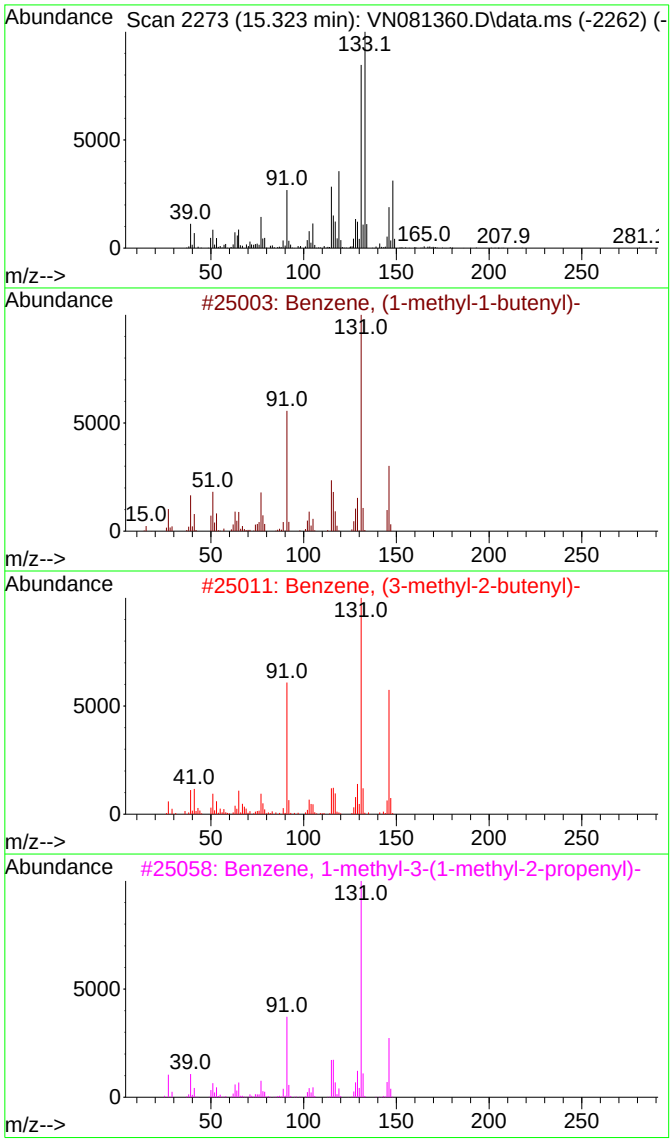
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

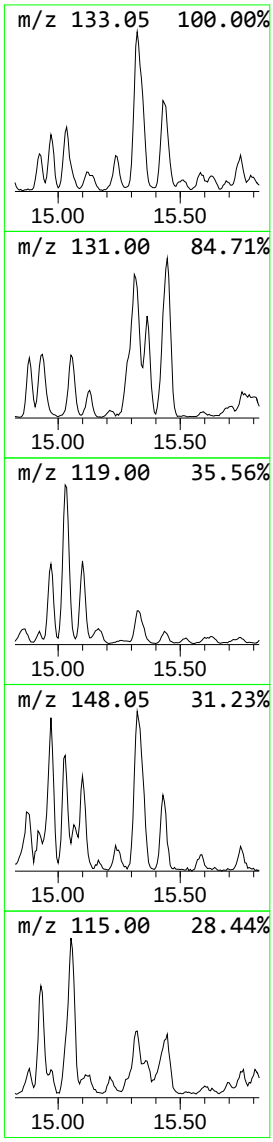
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, (1-methyl-1-butenyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.323	89.57 ug/l	2922950	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
3	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	50
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	50
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

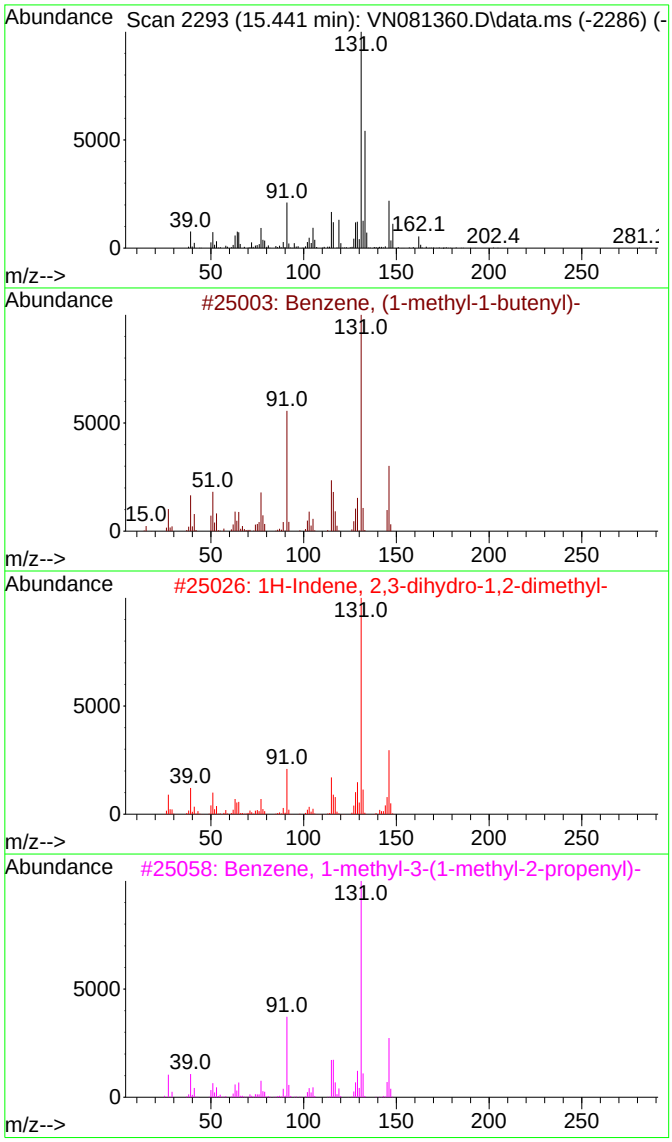
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.M
Quant Title : SW846 8260

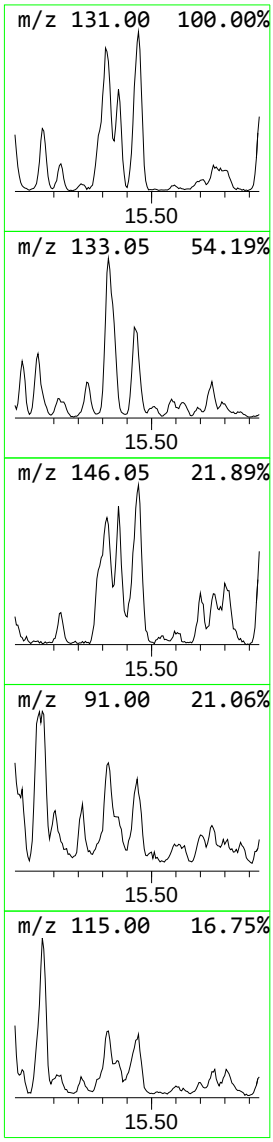
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.441	51.64 ug/l	1685120	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	91
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	83
3	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	64
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	64
5	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031124\
Data File : VN081360.D
Acq On : 11 Mar 2024 13:52
Operator : JC\MD
Sample : P1724-02
Misc : 1.02g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
10252

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030524W.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
unknown13.988	13.988	46.8	ug/l	1528050	4	13.794	1631680	50.0
o-Cymene	14.247	45.0	ug/l	1468400	4	13.794	1631680	50.0
Benzene, 2-ethy...	14.335	85.6	ug/l	2792010	4	13.794	1631680	50.0
1-Phenyl-1-butene	14.447	50.0	ug/l	1633380	4	13.794	1631680	50.0
Benzene, 1,2,4,...	14.659	65.9	ug/l	2150120	4	13.794	1631680	50.0
Benzene, 2-ethy...	14.694	65.8	ug/l	2146190	4	13.794	1631680	50.0
Benzene, 2-ethe...	14.929	55.1	ug/l	1799720	4	13.794	1631680	50.0
Benzene, 1-meth...	15.053	121.2	ug/l	3955030	4	13.794	1631680	50.0
Benzene, (1-met...	15.323	89.6	ug/l	2922950	4	13.794	1631680	50.0
1H-Indene, 2,3-...	15.441	51.6	ug/l	1685120	4	13.794	1631680	50.0