LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100324\
Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
Sample : P4275-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

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\82X100124W.M
Title : SW846 8260

Signal : TIC: VX043263.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.240	22	25	39	rBV	33381	49716	9.90%	0.978%
2	5.385	695	705	722	rBV	59322	174382	34.73%	3.431%
3	5.544	722	731	749	rVB2	88449	258809	51.54%	5.093%
4	5.952	789	798	812	rBV2	68303	183083	36.46%	3.603%
5	6.757	921	930	949	rBV	153262	373337	74.35%	7.346%
6	8.647	1234	1240	1257	rBV	317275	502167	100.00%	9.881%
7	10.055	1465	1471	1487	rBV	300309	423058	84.25%	8.325%
8	11.079	1634	1639	1652	rBV	233472	310312	61.79%	6.106%
9	11.616	1720	1727	1733	rBV5	3248	5550	1.11%	0.109%
10	12.018	1788	1793	1805	rBV	267417	358212	71.33%	7.049%
11	12.274	1832	1835	1838	rBV2	5013	5838	1.16%	0.115%
12	12.317	1838	1842	1848	rVB2	20847	32501	6.47%	0.640%
13	12.421	1848	1859	1863	rBV6	5565	12497	2.49%	0.246%
14	12.463	1863	1866	1873	rVB4	3533	5610	1.12%	0.110%
15	12.610	1884	1890	1895	rVB	24079	31943	6.36%	0.629%
16	12.725	1903	1909	1913	rBV2	19138	33771	6.73%	0.665%
17	12.799	1917	1921	1926	rVV2	14353	29164	5.81%	0.574%
18	12.841	1926	1928	1932	rVB2	7264	8753	1.74%	0.172%
19	12.920	1938	1941	1945	rVB	14226	16464	3.28%	0.324%
20	12.963	1945	1948	1950	rBV2	4503	5652	1.13%	0.111%
21	13.042	1954	1961	1964	rBV2	31404	61014	12.15%	1.201%
22	13.067	1964	1965	1968	rVV	14686	14044	2.80%	0.276%
23	13.091	1968	1969	1973	rVB	10863	9231	1.84%	0.182%
24	13.140	1973	1977	1983	rVB3	20363	29339	5.84%	0.577%
25	13.207	1983	1988	1991	rBV4	7754	11111	2.21%	0.219%
26	13.262	1991	1997	2000	rBV3	22429	36143	7.20%	0.711%
27	13.298	2000	2003	2007	rVV4	12034	24037	4.79%	0.473%
28	13.341	2007	2010	2012	rVV4	12294	16877	3.36%	0.332%
29	13.372	2012	2015	2021	rVB3	12846	18480	3.68%	0.364%
30	13.433	2021	2025	2032	rVB3	19009	32919	6.56%	0.648%
31	13.500	2032	2036	2039	rBV4	8073	13386	2.67%	0.263%
32	13.548	2039	2044	2046	rBV	42289	55202	10.99%	1.086%
33	13.579	2046	2049	2055	rVV3	59607	102057	20.32%	2.008%
34	13.640	2055	2059	2060	rVV	27566	32032	6.38%	0.630%
35	13.664	2060	2063	2068	rVB3	36976	64533	12.85%	1.270%
36	13.713	2068	2071	2074	rBV	4664	5963	1.19%	0.117%

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ClientSampleId :
TRE-1768

Integration Parameters: RTEINT.P
Integrator: RTE
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Title : SW846 8260

37	13.750	2074	2077	2081	rVB	11707	14679	2.92%	0.289%
38	13.792	2081	2084	2088	rBV3	14153	17943	3.57%	0.353%
39	13.853	2088	2094	2099	rVB4	33776	64364	12.82%	1.266%
40	13.926	2099	2106	2110	rBV2	48762	79024	15.74%	1.555%
41	13.969	2110	2113	2116	rBV2	14407	17746	3.53%	0.349%
42	14.006	2116	2119	2121	rBV	8028	7133	1.42%	0.140%
43	14.036	2121	2124	2127	rVB2	17865	17056	3.40%	0.336%
44	14.073	2127	2130	2134	rVB3	15452	19095	3.80%	0.376%
45	14.164	2137	2145	2149	rBV4	14969	35571	7.08%	0.700%
46	14.213	2149	2153	2158	rBV4	67263	104224	20.75%	2.051%
47	14.323	2167	2171	2176	rBV2	44487	69645	13.87%	1.370%
48	14.371	2176	2179	2183	rVV	87158	102586	20.43%	2.019%
49	14.414	2183	2186	2192	rVB2	23577	37851	7.54%	0.745%
50	14.481	2192	2197	2202	rBV2	61242	106161	21.14%	2.089%
51	14.524	2202	2204	2209	rVB4	7968	9088	1.81%	0.179%
52	14.579	2210	2213	2215	rBV3	9185	11554	2.30%	0.227%
53	14.615	2215	2219	2225	rVV2	70863	122178	24.33%	2.404%
54	14.670	2225	2228	2233	rVB	68239	87055	17.34%	1.713%
55	14.725	2233	2237	2239	rBV	23690	30332	6.04%	0.597%
56	14.755	2239	2242	2245	rVB	17737	17164	3.42%	0.338%
57	14.816	2245	2252	2255	rBV2	58757	98431	19.60%	1.937%
58	14.841	2255	2256	2261	rVB3	38213	40392	8.04%	0.795%
59	14.902	2261	2266	2272	rBV2	46643	88748	17.67%	1.746%
60	14.981	2277	2279	2283	rVB4	15591	21701	4.32%	0.427%
61	15.048	2285	2290	2294	rVB3	19714	29222	5.82%	0.575%
62	15.103	2295	2299	2302	rBV5	10674	13898	2.77%	0.273%
63	15.146	2302	2306	2308	rBV4	8840	14135	2.81%	0.278%
64	15.182	2308	2312	2319	rVV	93181	166559	33.17%	3.277%
65	15.243	2319	2322	2325	rVV4	8168	9775	1.95%	0.192%
66	15.274	2325	2327	2330	rVV3	7456	7799	1.55%	0.153%
67	15.316	2330	2334	2339	rVB5	9467	17184	3.42%	0.338%
68	15.371	2339	2343	2352	rVB5	18651	37840	7.54%	0.745%
69	15.444	2352	2355	2357	rBV4	5224	6001	1.20%	0.118%
70	15.481	2357	2361	2365	rBV3	27643	47459	9.45%	0.934%
71	15.524	2365	2368	2372	rVV3	21417	31148	6.20%	0.613%
72	15.566	2372	2375	2377	rVV4	7936	9439	1.88%	0.186%
73	15.609	2379	2382	2386	rVB3	17365	24162	4.81%	0.475%
74	15.755	2396	2406	2408	rBV7	9796	25759	5.13%	0.507%
75	15.847	2416	2421	2424	rBV5	10932	20297	4.04%	0.399%
76	15.999	2441	2446	2447	rBV5	3629	6432	1.28%	0.127%

LSC Area Percent Report

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ClientSampleId :
TRE-1768

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Title : SW846 8260

77	16.121	2462	2466	2472	rVB8	4464	9242	1.84%	0.182%
78	16.286	2486	2493	2500	rVB8	6181	16686	3.32%	0.328%
79	16.371	2504	2507	2512	rVB7	4550	6552	1.30%	0.129%
80	16.511	2527	2530	2533	rBV3	4102	6447	1.28%	0.127%
81	16.712	2559	2563	2569	rVB7	4365	9130	1.82%	0.180%

Sum of corrected areas: 5082074

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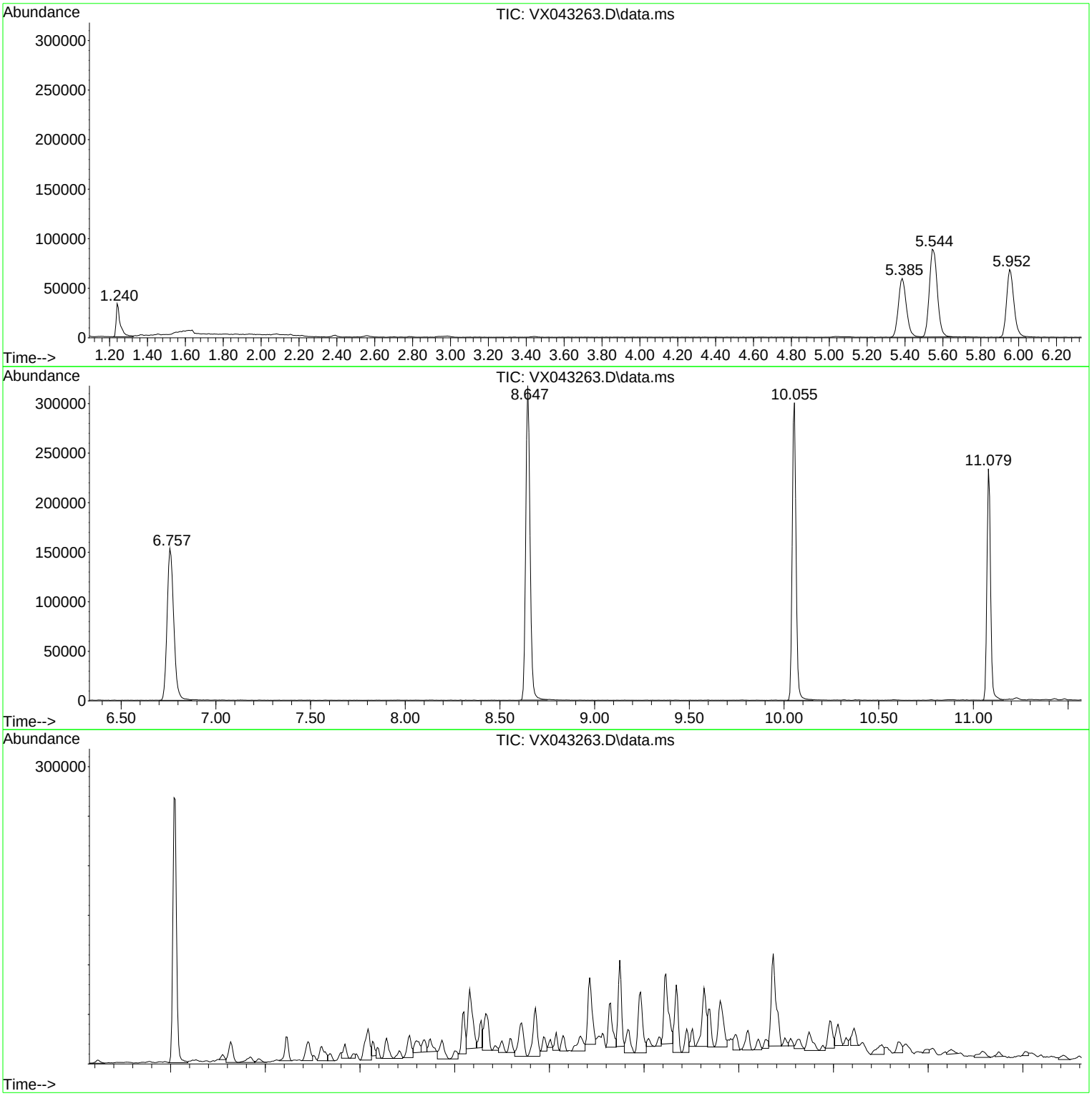
ClientSampleId :
TRE-1768

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

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



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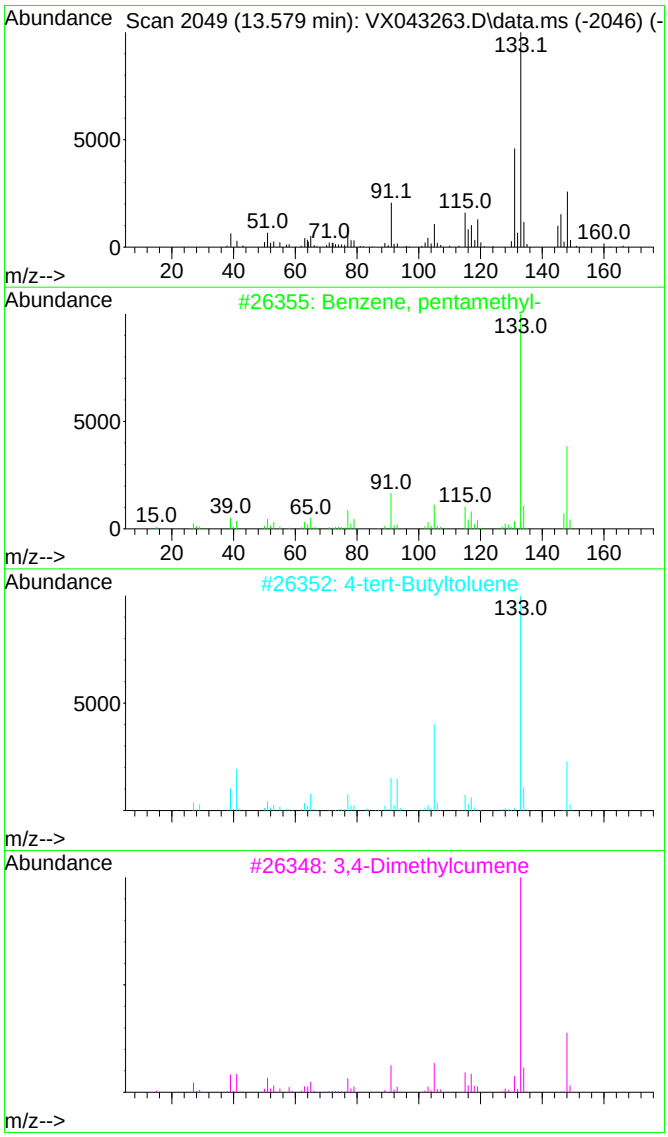
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

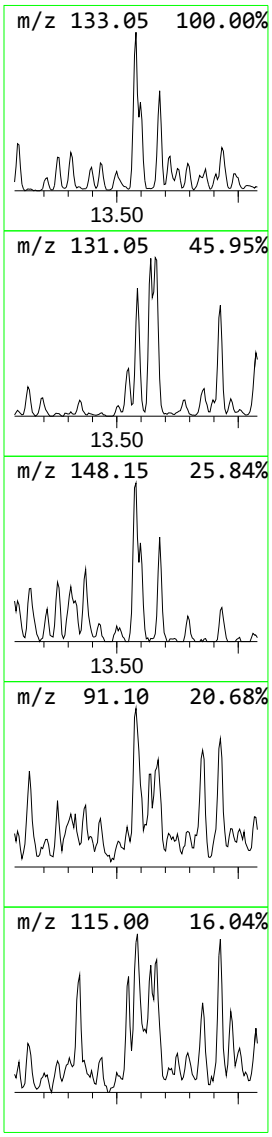
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, pentamethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	14.25 ug/l	102057	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	58
2	4-tert-Butyltoluene	148	C11H16	000098-51-1	53
3	3,4-Dimethylcumene	148	C11H16	1000370-34-1	52
4	Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	52
5	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	52





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ClientSampleId :
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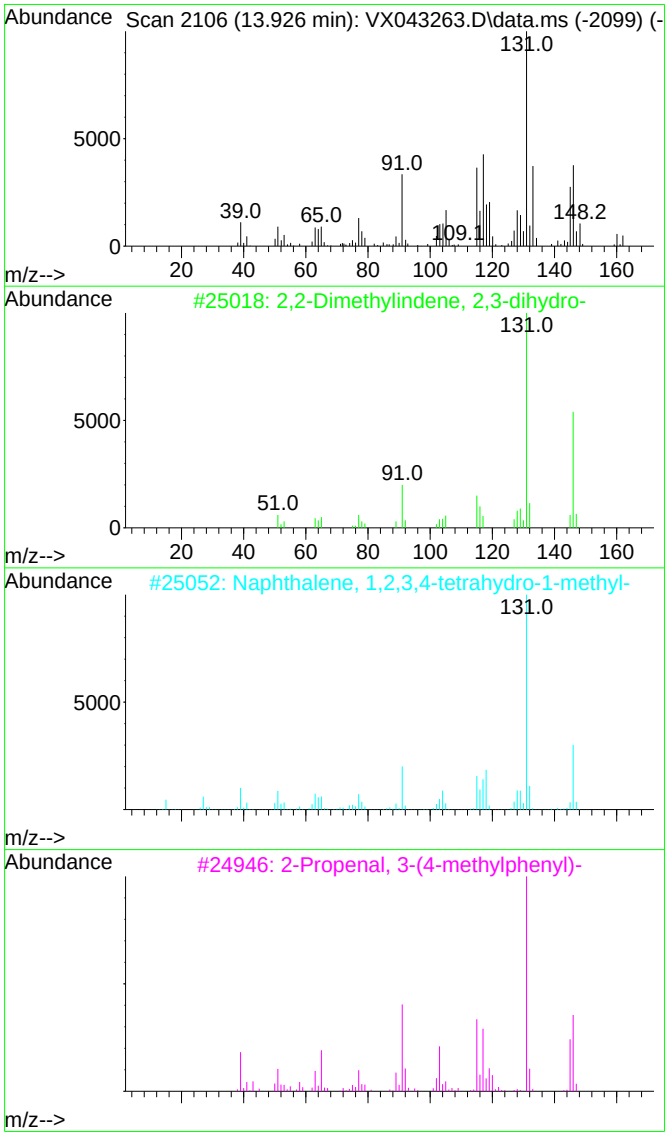
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

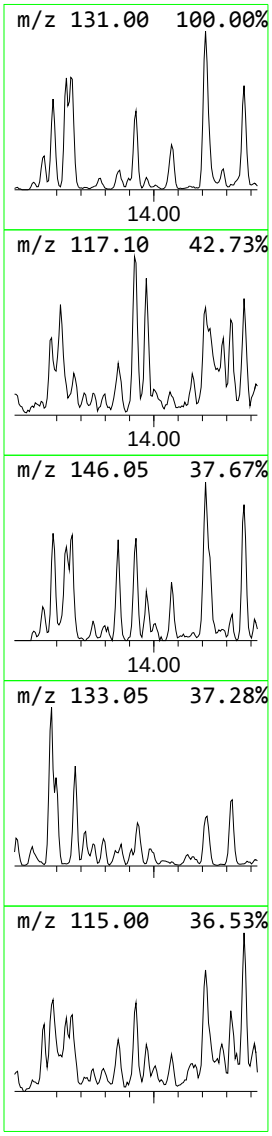
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2,2-Dimethylindene, 2,3-dih... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	11.03 ug/l	79024	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60
3	2-Propenal, 3-(4-methylphenyl)-	146	C10H10O	001504-75-2	58
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	55
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	160	C12H16	078920-29-3	53





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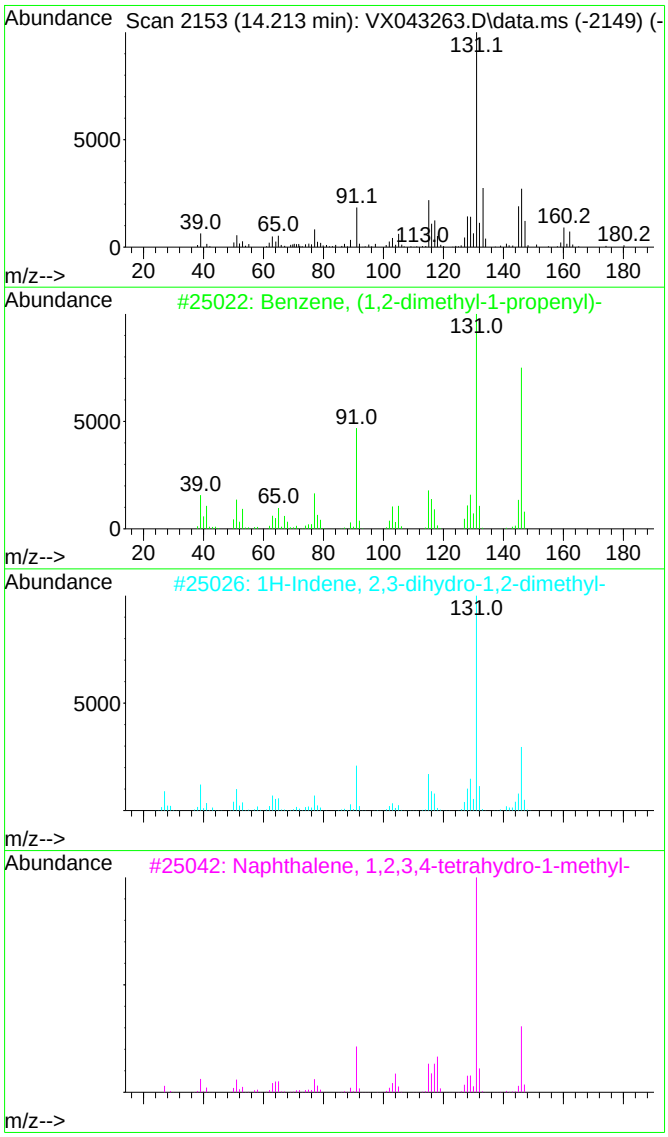
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

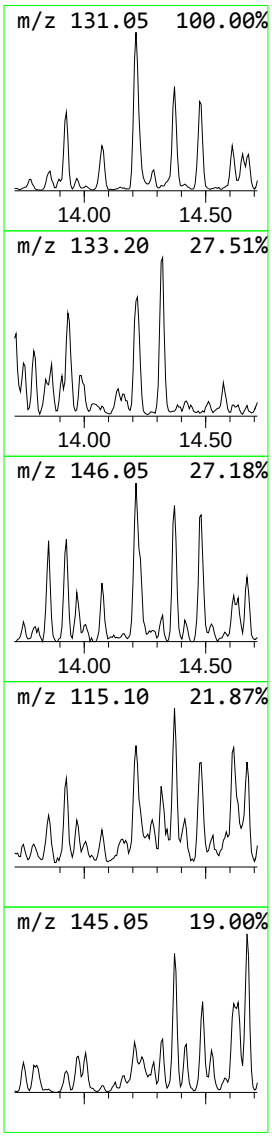
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	14.55 ug/l	104224	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76





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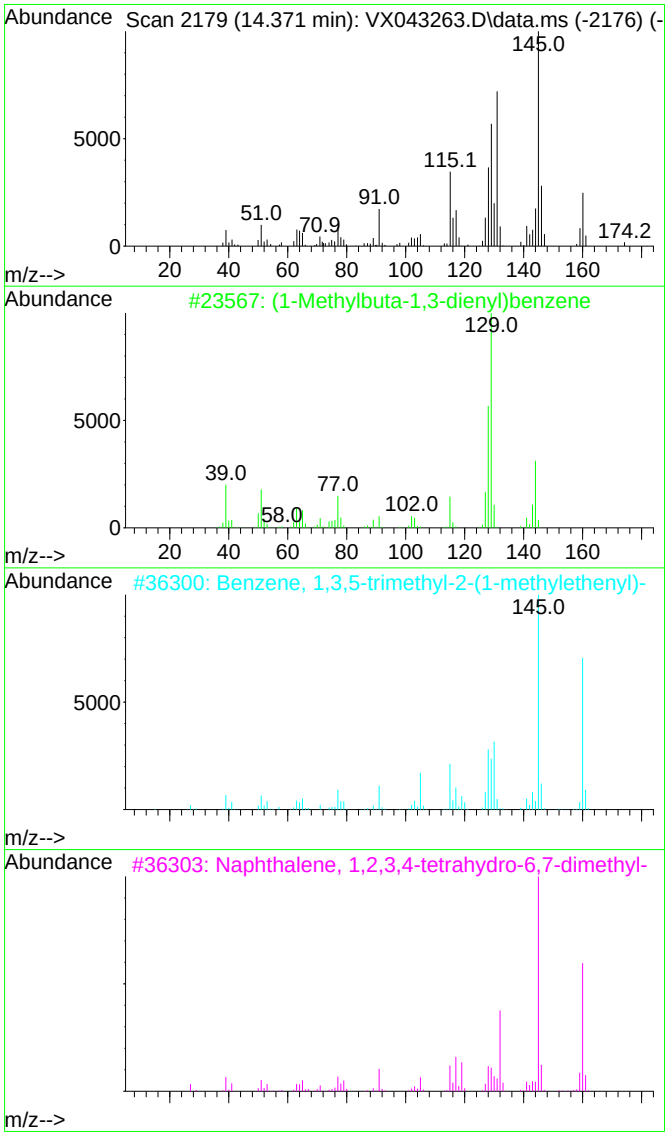
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

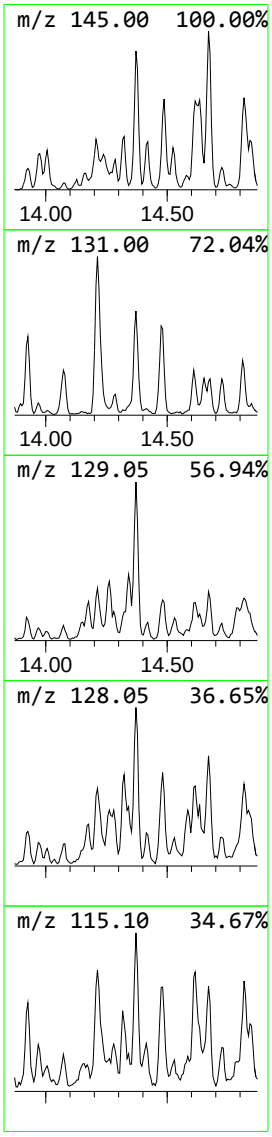
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (1-Methylbuta-1,3-dienyl)be... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.371	14.32 ug/l	102586	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1-Methylbuta-1,3-dienyl)benzene	144	C11H12	054758-36-0	60
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	53
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	46
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	46
5	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	001198-20-5	45





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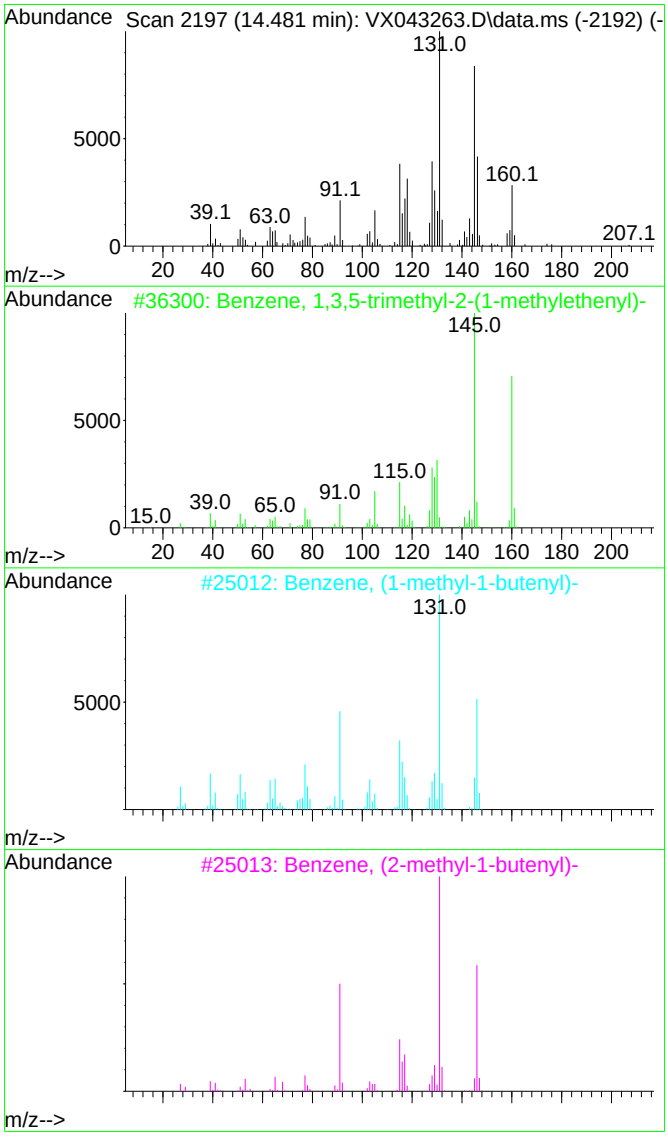
Quant Title : SW846 8260

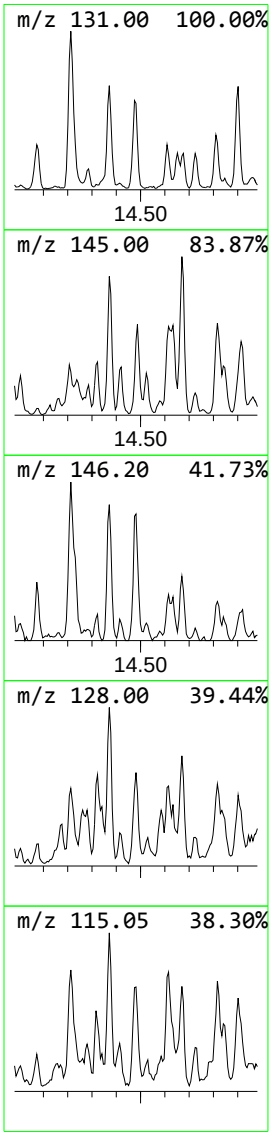
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	14.82 ug/l	106161	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	60
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	53
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	46
4	Ethanone, 1-(2,3-dihydro-1H-inde...	160	C11H12O	004228-10-8	46
5	Ethanone, 1-[4-(1-methylethenyl)...	160	C11H12O	005359-04-6	42





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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
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TIC Library : C:\Database\NIST20.L
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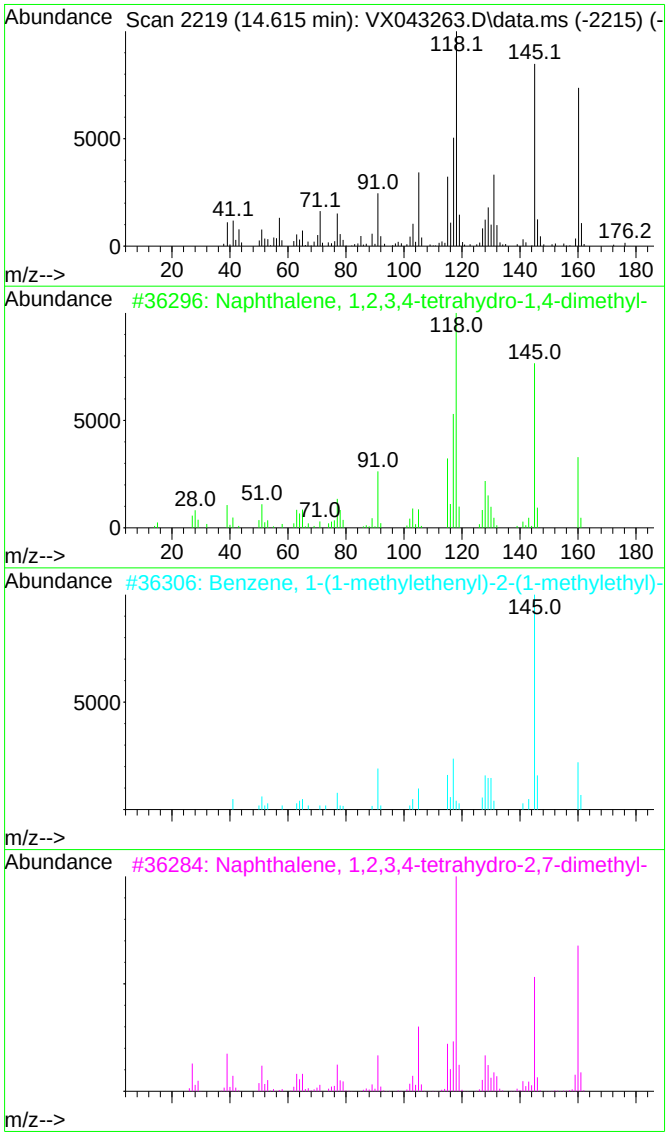
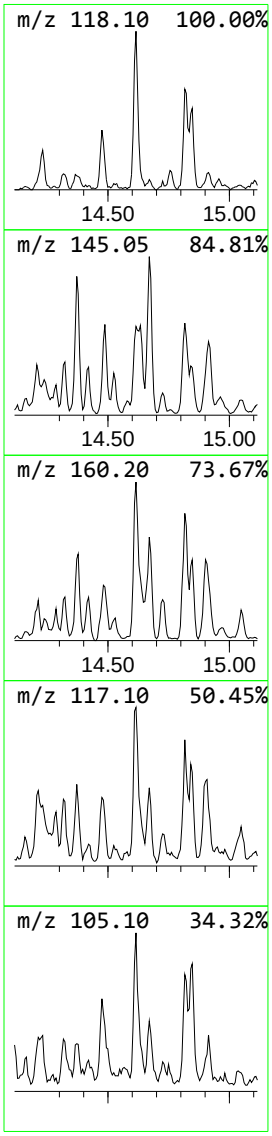
****
Peak Number 6 Benzene, 1-(1-methylethenyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	17.05 ug/l	122178	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	97
2	Benzene, 1-(1-methylethenyl)-2-(...	160	C12H16	005557-93-7	83
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	70
4	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	55
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	55





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Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
Sample : P4275-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

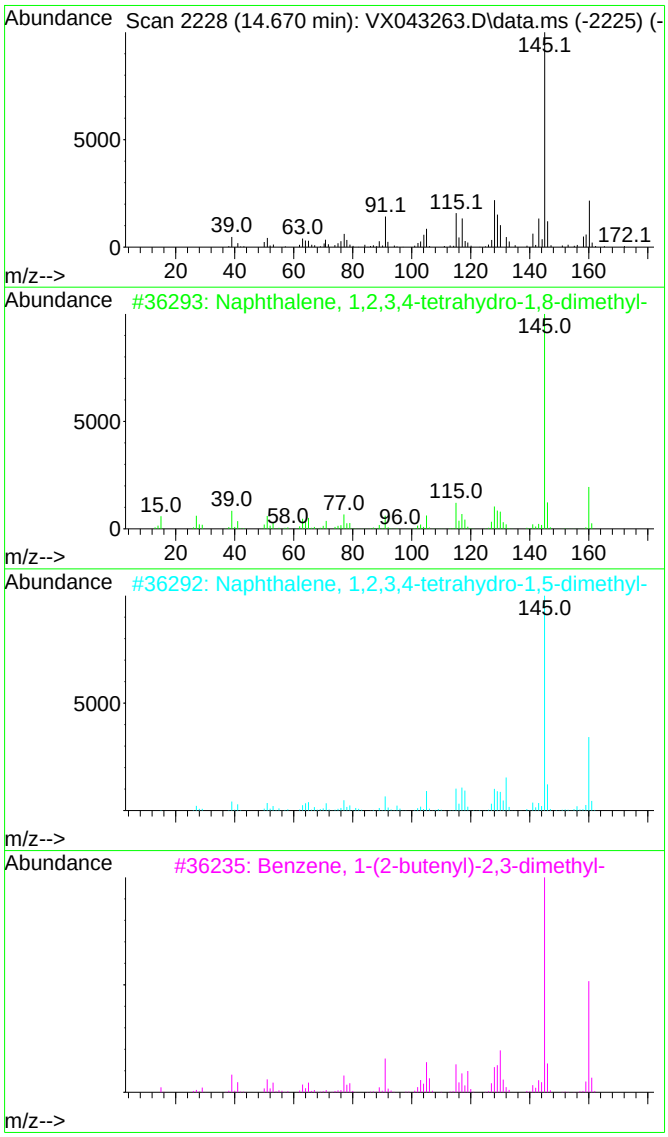
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

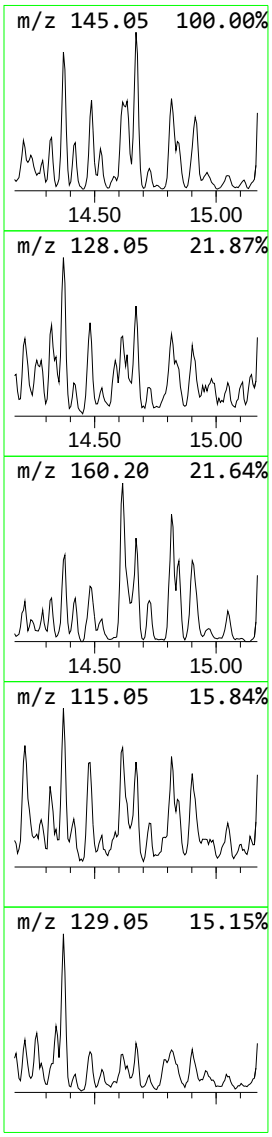
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.670	12.15 ug/l	87055	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	83
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	80
3	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	72
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001985-59-7	72
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100324\
Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
Sample : P4275-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

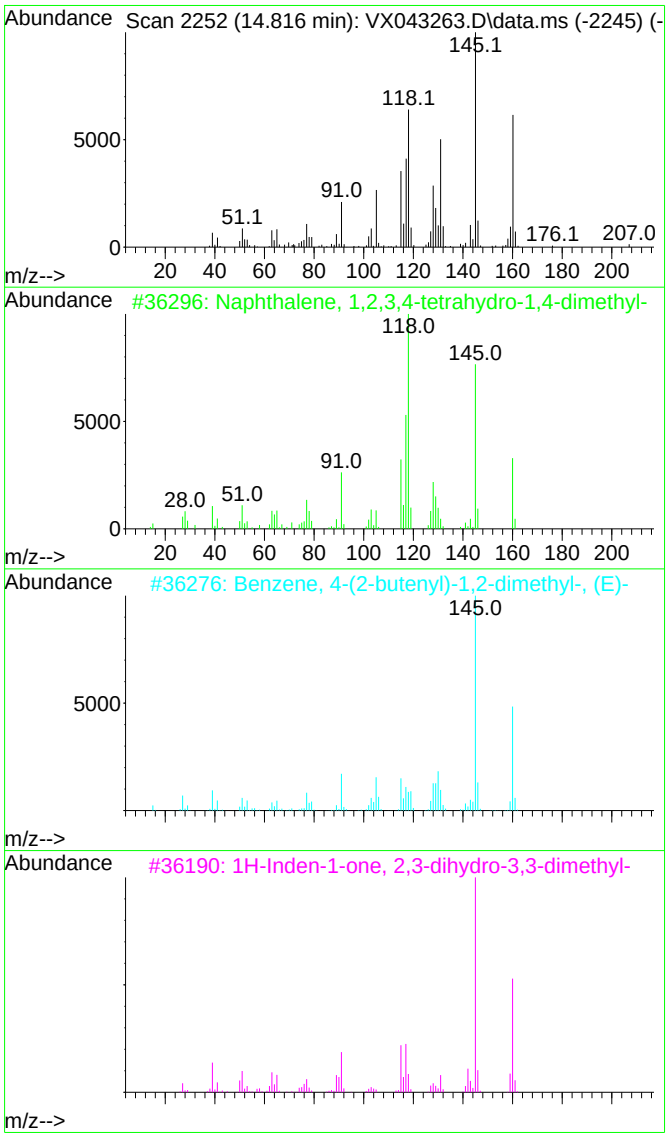
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

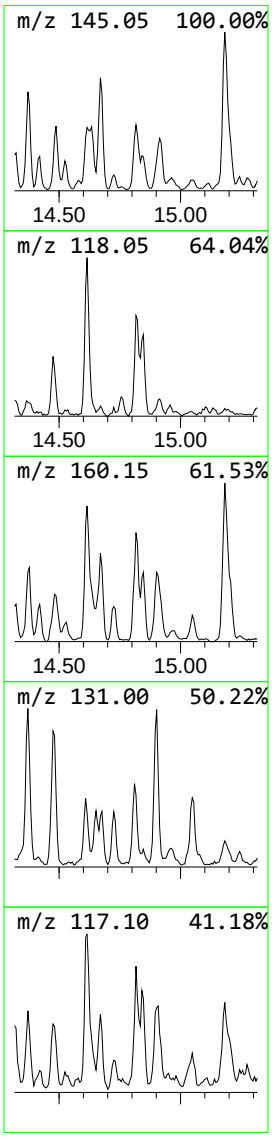
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.816	13.74 ug/l	98431	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	94
2	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	68
3	1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	62
4	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	60
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100324\
Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
Sample : P4275-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

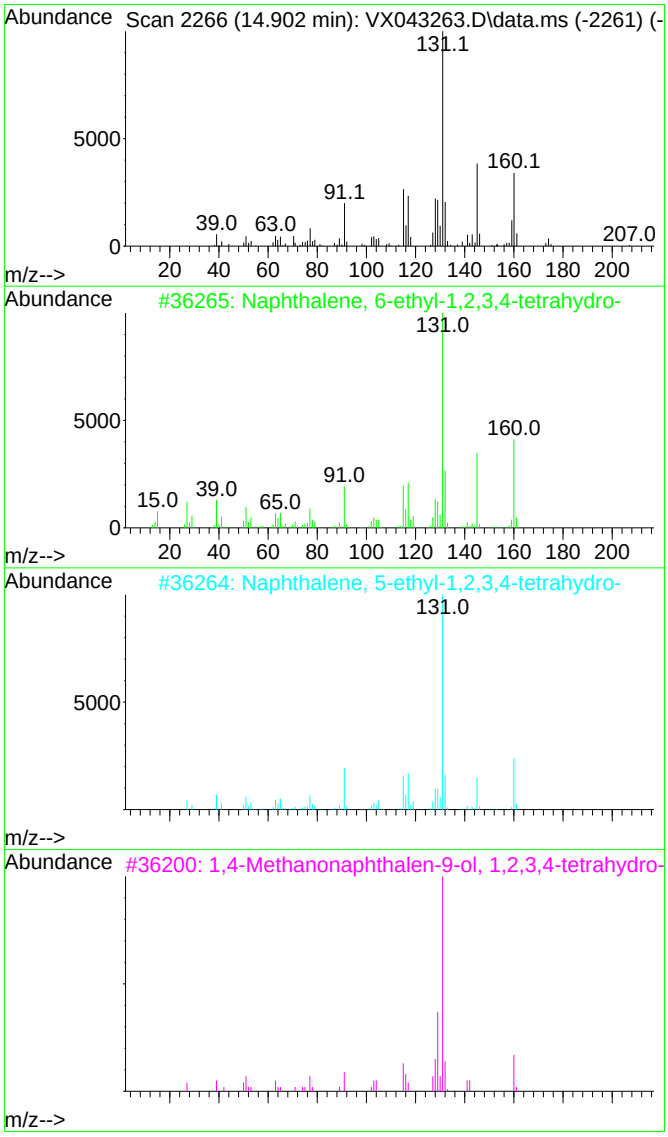
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

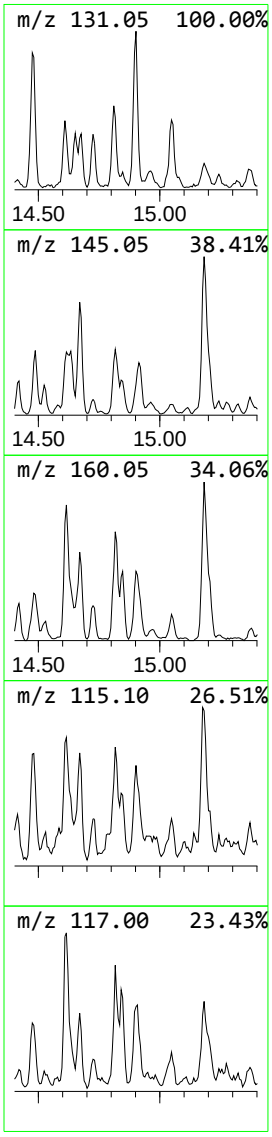
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.902	12.39 ug/l	88748	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	87
2	Naphthalene, 5-ethyl-1,2,3,4-tet...	160	C12H16	042775-75-7	68
3	1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	055255-94-2	43
4	Benzene, (3-chloro-1-methyl-1-pr...	166	C10H11Cl	1000327-43-7	38
5	1,2,3,4-Tetrahydro-1-naphthalene...	236	C13H20SSi	1000470-19-9	38





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Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

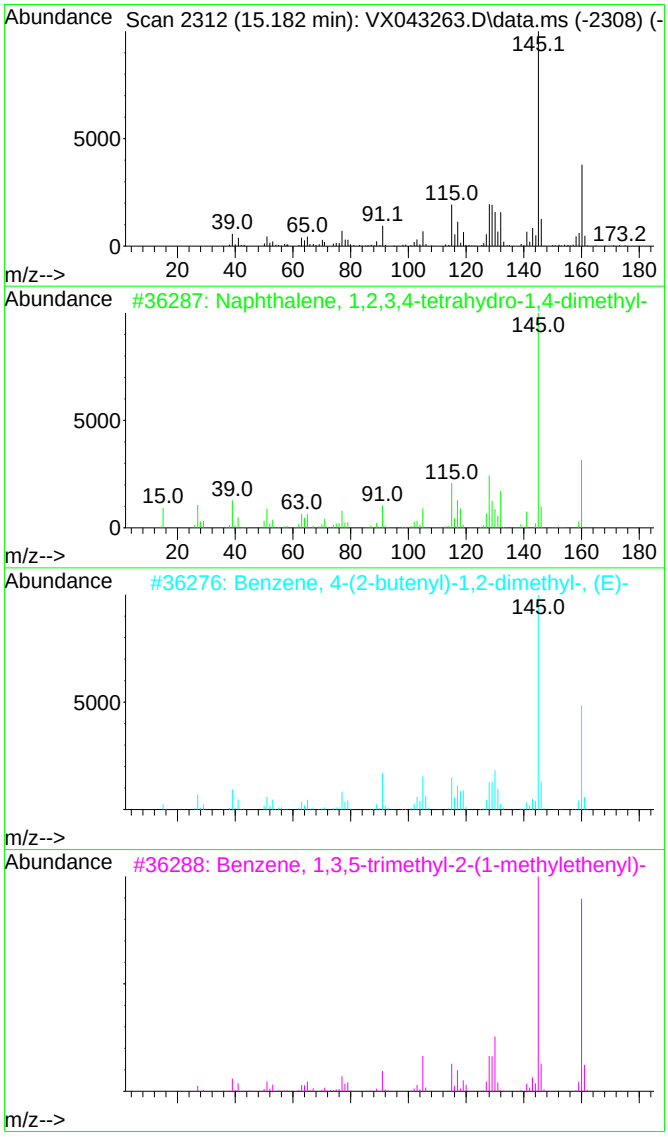
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
Quant Title : SW846 8260

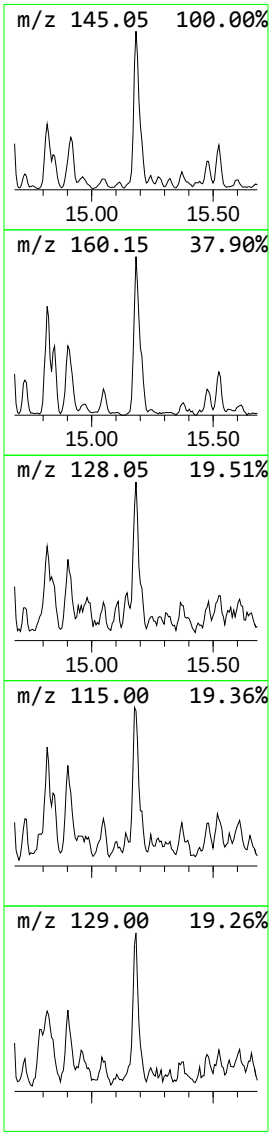
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-(2-butenyl)-2,3-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	23.25 ug/l	166559	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	93
2	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	91
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	90
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	89
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100324\
Data File : VX043263.D
Acq On : 03 Oct 2024 15:35
Operator : JC/MD
Sample : P4275-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1768

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.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
Benzene, pentam...	13.579	14.3	ug/l	102057	4	12.024	358212	50.0
2,2-Dimethylind...	13.926	11.0	ug/l	79024	4	12.024	358212	50.0
Benzene, (1,2-d...	14.213	14.6	ug/l	104224	4	12.024	358212	50.0
(1-Methylbuta-1...	14.371	14.3	ug/l	102586	4	12.024	358212	50.0
Benzene, 1,3,5-...	14.481	14.8	ug/l	106161	4	12.024	358212	50.0
Benzene, 1-(1-m...	14.615	17.1	ug/l	122178	4	12.024	358212	50.0
Naphthalene, 1,...	14.670	12.2	ug/l	87055	4	12.024	358212	50.0
Benzene, 4-(2-b...	14.816	13.7	ug/l	98431	4	12.024	358212	50.0
Naphthalene, 6-...	14.902	12.4	ug/l	88748	4	12.024	358212	50.0
Benzene, 1-(2-b...	15.182	23.3	ug/l	166559	4	12.024	358212	50.0