

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
 Data File : VV024340.D
 Acq On : 12 Jan 2022 19:10
 Operator : SY/MD
 Sample : N1084-09 20X
 Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLN2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % 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_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

Signal : TIC: VV024340.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.310	76	84	95	rBV	128451	134730	2.59%	0.280%
2	1.571	157	165	179	rBV	107258	123958	2.38%	0.258%
3	2.108	321	332	359	rVB	245808	357791	6.87%	0.745%
4	2.503	450	455	458	rBV3	1518	1599	0.03%	0.003%
5	2.767	527	537	546	rBV5	1988	3872	0.07%	0.008%
6	3.047	614	624	639	rVB3	3853	8061	0.15%	0.017%
7	3.272	690	694	701	rVB2	262	418	0.01%	0.001%
8	3.767	836	848	861	rBV8	2576	6003	0.12%	0.012%
9	3.889	876	886	917	rBV	55590	171063	3.28%	0.356%
10	4.352	1012	1030	1049	rBV	155267	379638	7.29%	0.790%
11	4.516	1075	1081	1085	rBV3	345	331	0.01%	0.001%
12	4.558	1089	1094	1097	rVB3	499	437	0.01%	0.001%
13	4.674	1120	1130	1143	rVB5	3486	8135	0.16%	0.017%
14	4.770	1152	1160	1162	rBV4	1240	1571	0.03%	0.003%
15	4.912	1192	1204	1218	rBV4	5295	12913	0.25%	0.027%
16	5.047	1229	1246	1281	rBV2	357350	952338	18.29%	1.983%
17	5.323	1323	1332	1338	rBV7	1255	2093	0.04%	0.004%
18	5.394	1348	1354	1358	rBV	334	359	0.01%	0.001%
19	5.490	1375	1384	1404	rBV4	10984	23514	0.45%	0.049%
20	5.616	1410	1423	1448	rBV	290751	579611	11.13%	1.207%
21	5.911	1504	1515	1528	rVB4	8353	16967	0.33%	0.035%
22	5.956	1528	1529	1533	rVB2	570	310	0.01%	0.001%
23	6.069	1551	1564	1580	rBV	197853	408238	7.84%	0.850%
24	6.256	1619	1622	1625	rBV3	1108	724	0.01%	0.002%
25	6.368	1651	1657	1661	rBV4	588	533	0.01%	0.001%
26	6.465	1683	1687	1690	rBV2	595	505	0.01%	0.001%
27	6.635	1733	1740	1741	rBV4	976	814	0.02%	0.002%
28	6.651	1743	1745	1749	rVB2	585	474	0.01%	0.001%
29	6.776	1779	1784	1785	rBV3	884	623	0.01%	0.001%
30	6.834	1798	1802	1809	rVB5	766	946	0.02%	0.002%
31	6.918	1818	1828	1837	rBV5	4254	7488	0.14%	0.016%
32	6.985	1838	1849	1869	rBV	118662	210516	4.04%	0.438%
33	7.265	1923	1936	1940	rBV4	6377	10859	0.21%	0.023%
34	7.317	1941	1952	1967	rVV	452256	777660	14.93%	1.619%
35	7.391	1968	1975	1988	rVB2	35067	61293	1.18%	0.128%
36	7.567	2019	2030	2038	rBV3	13416	21116	0.41%	0.044%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

37	7.622	2038	2047	2071	rVB	82866	152578	2.93%	0.318%
38	7.789	2091	2099	2110	rVB7	3074	5029	0.10%	0.010%
39	7.908	2126	2136	2139	rBV5	1496	2235	0.04%	0.005%
40	7.976	2150	2157	2169	rVB8	3456	5732	0.11%	0.012%
41	8.088	2174	2192	2220	rBV	230252	423354	8.13%	0.881%
42	8.204	2224	2228	2235	rBV5	5233	8767	0.17%	0.018%
43	8.288	2246	2254	2264	rVB2	17122	27906	0.54%	0.058%
44	8.352	2264	2274	2283	rBV2	24098	39166	0.75%	0.082%
45	8.509	2312	2323	2332	rBV5	4657	7691	0.15%	0.016%
46	8.571	2332	2342	2344	rBV3	17775	23325	0.45%	0.049%
47	8.667	2364	2372	2384	rVB3	54227	92449	1.78%	0.192%
48	8.789	2399	2410	2419	rBV	52707	87925	1.69%	0.183%
49	8.850	2419	2429	2445	rBV	451686	718182	13.79%	1.495%
50	9.011	2465	2479	2493	rBV	890644	1386943	26.63%	2.887%
51	9.136	2506	2518	2531	rBV	3287226	5207533	100.00%	10.841%
52	9.204	2531	2539	2563	rVB2	304218	502441	9.65%	1.046%
53	9.316	2563	2574	2589	rBV4	22050	39203	0.75%	0.082%
54	9.410	2595	2603	2606	rBV3	10564	14072	0.27%	0.029%
55	9.474	2617	2623	2633	rVB4	94613	139252	2.67%	0.290%
56	9.542	2634	2644	2665	rVB	842154	1345039	25.83%	2.800%
57	9.673	2678	2685	2688	rBV3	43441	56093	1.08%	0.117%
58	9.709	2689	2696	2706	rVB2	225939	346295	6.65%	0.721%
59	9.799	2716	2724	2733	rBV4	179831	288880	5.55%	0.601%
60	10.017	2785	2792	2800	rBV3	97100	140992	2.71%	0.294%
61	10.217	2845	2854	2872	rVB2	507619	1011480	19.42%	2.106%
62	10.352	2887	2896	2904	rBV	397312	607309	11.66%	1.264%
63	10.445	2916	2925	2940	rBV	1455075	3000183	57.61%	6.246%
64	10.538	2948	2954	2960	rBV	752276	977539	18.77%	2.035%
65	10.577	2961	2966	2987	rVB	705774	1066193	20.47%	2.220%
66	10.734	3006	3015	3035	rVB	609931	1065952	20.47%	2.219%
67	10.879	3052	3060	3062	rBV	313097	380256	7.30%	0.792%
68	10.911	3063	3070	3082	rVB	3079524	4560963	87.58%	9.495%
69	11.033	3101	3108	3112	rBV	105504	159293	3.06%	0.332%
70	11.152	3137	3145	3153	rBV5	332319	640398	12.30%	1.333%
71	11.246	3166	3174	3183	rBV2	616496	931419	17.89%	1.939%
72	11.336	3194	3202	3209	rBV	914170	1280662	24.59%	2.666%
73	11.461	3236	3241	3250	rVB2	181412	248818	4.78%	0.518%
74	11.525	3250	3261	3271	rBV	2975295	4728672	90.80%	9.845%
75	11.628	3283	3293	3310	rVB6	946464	2249325	43.19%	4.683%
76	11.911	3373	3381	3384	rBV	243274	327390	6.29%	0.682%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Title : VOC Analysis

77	12.014	3407	3413	3436	rVB	360405	639559	12.28%	1.331%
78	12.117	3438	3445	3449	rBV2	95095	140970	2.71%	0.293%
79	12.255	3477	3488	3497	rBV6	123779	246701	4.74%	0.514%
80	12.368	3514	3523	3530	rBV7	150340	284728	5.47%	0.593%
81	12.413	3532	3537	3544	rVB	191848	252087	4.84%	0.525%
82	12.464	3544	3553	3563	rBV	439699	688778	13.23%	1.434%
83	12.725	3628	3634	3647	rVB2	131455	194836	3.74%	0.406%
84	12.795	3649	3656	3663	rBV2	41205	57566	1.11%	0.120%
85	12.847	3666	3672	3673	rBV4	29006	29890	0.57%	0.062%
86	12.882	3676	3683	3697	rVB3	223428	373122	7.17%	0.777%
87	13.001	3714	3720	3725	rBV2	37232	47156	0.91%	0.098%
88	13.046	3727	3734	3744	rVV6	72219	126367	2.43%	0.263%
89	13.268	3794	3803	3809	rBV4	49777	84381	1.62%	0.176%
90	13.500	3863	3875	3901	rVB	1310465	2086907	40.07%	4.345%
91	13.622	3904	3913	3933	rVB	81356	144856	2.78%	0.302%
92	13.905	3995	4001	4015	rVB5	26058	48160	0.92%	0.100%
93	14.091	4050	4059	4073	rBV5	24063	54865	1.05%	0.114%
94	14.574	4200	4209	4215	rBV3	21952	37910	0.73%	0.079%
95	14.628	4215	4226	4253	rVB	1865614	2836324	54.47%	5.905%
96	14.815	4273	4284	4300	rVB	556524	846383	16.25%	1.762%
97	15.361	4448	4454	4466	rVB	53054	78422	1.51%	0.163%
98	15.551	4507	4513	4520	rBV3	24019	34874	0.67%	0.073%
99	15.667	4540	4549	4563	rBV3	38803	75881	1.46%	0.158%
100	15.811	4587	4594	4600	rBV2	26806	38424	0.74%	0.080%

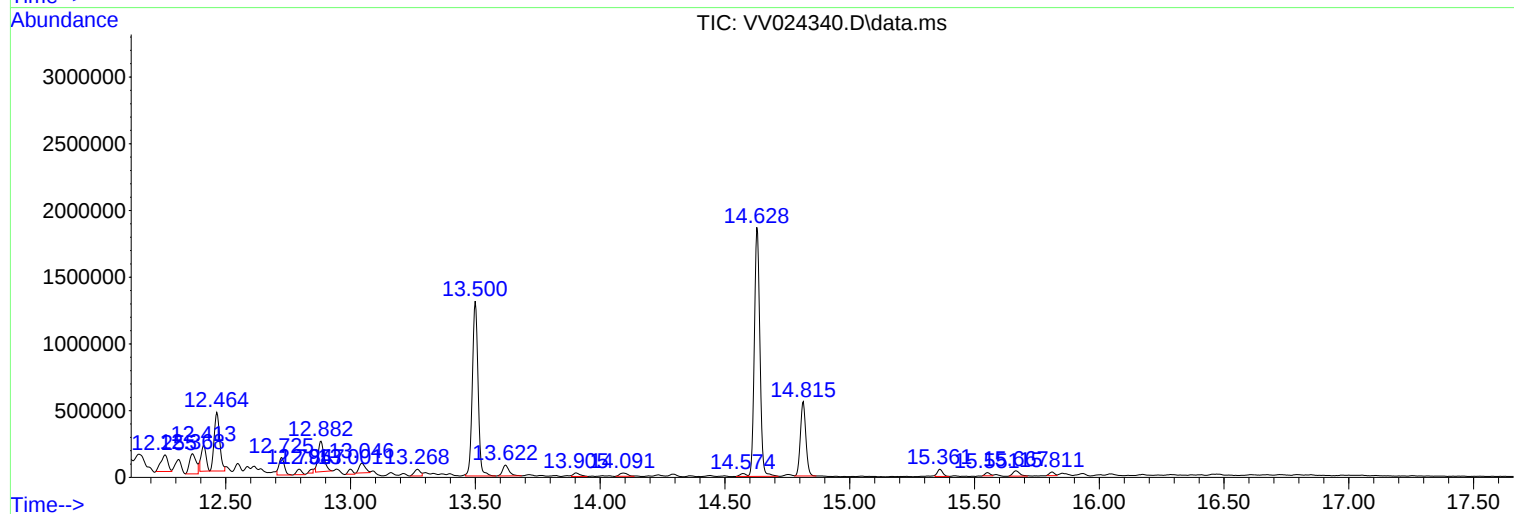
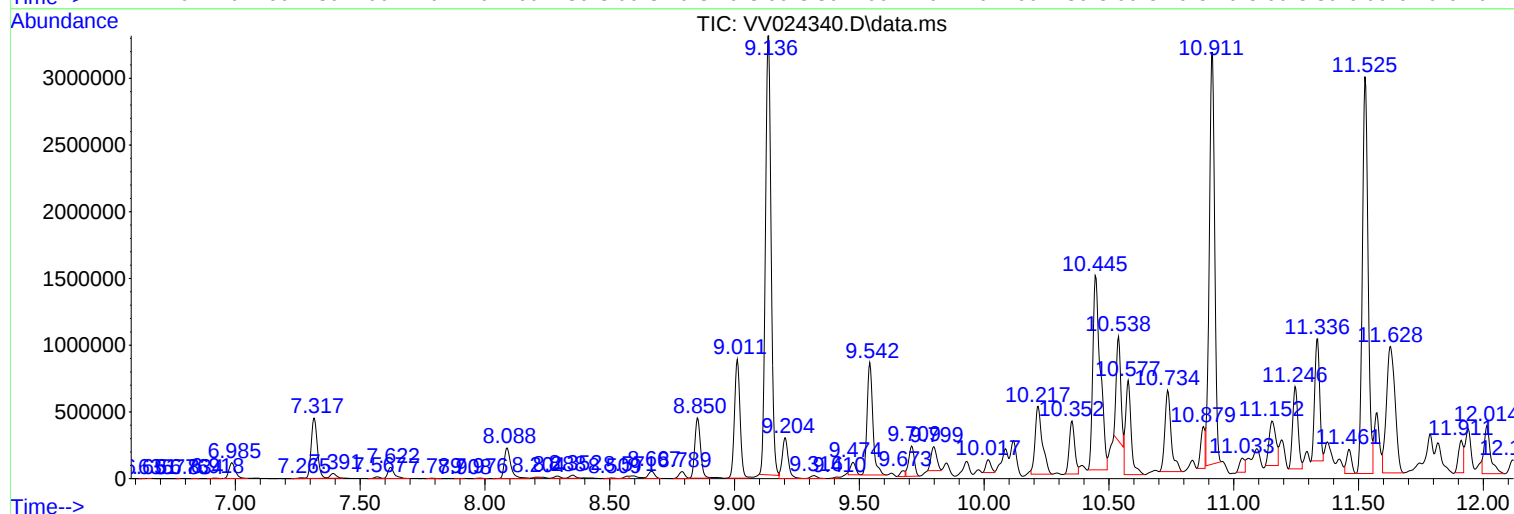
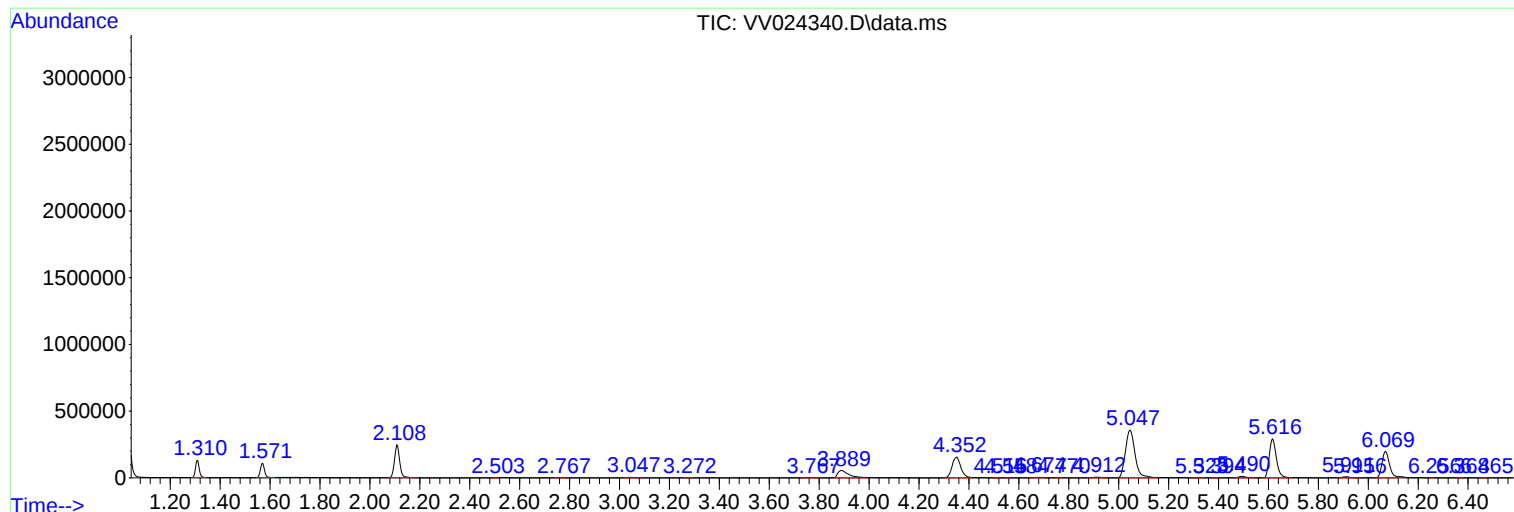
Sum of corrected areas: 48033582

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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
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Instrument :
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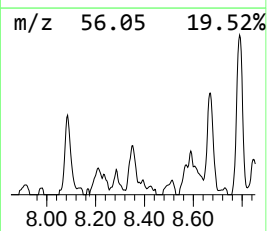
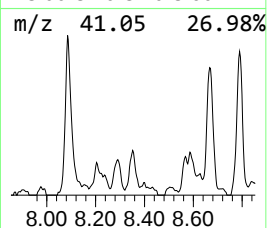
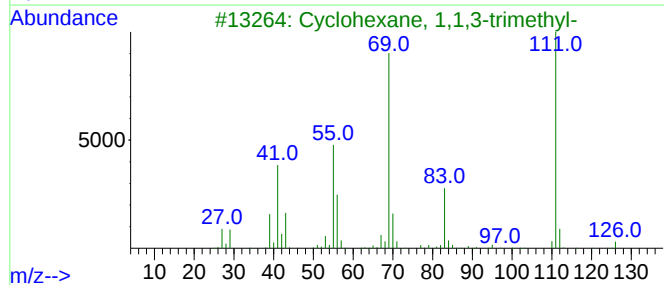
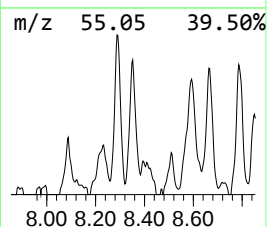
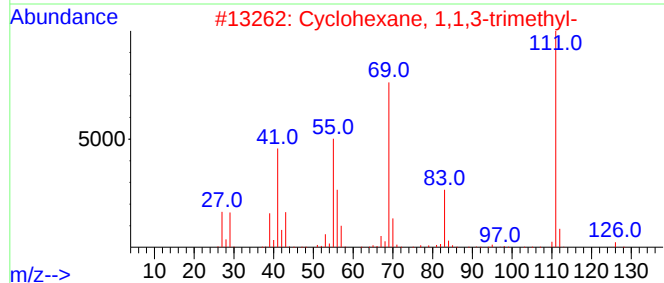
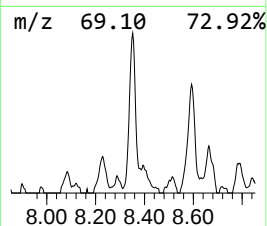
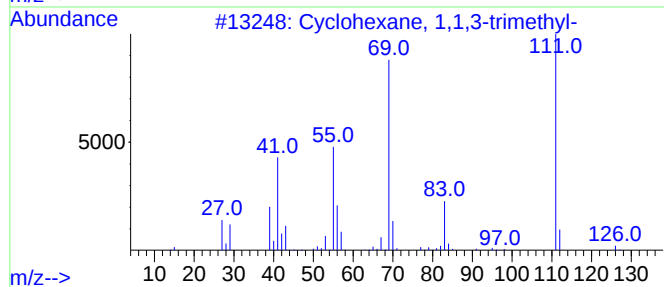
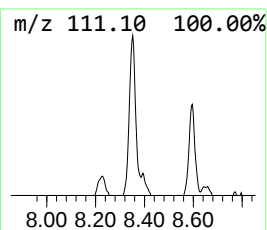
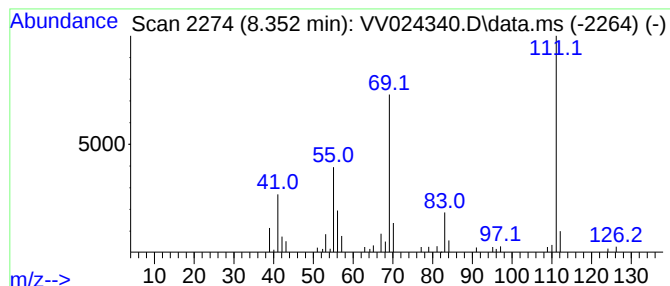
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic8.352 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.352	2.73 ug/L	39166	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	94
2			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	91
3			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	91
4			1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	80
5			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	70



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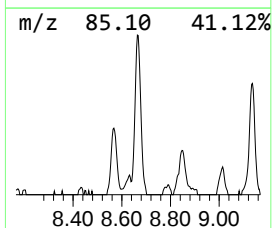
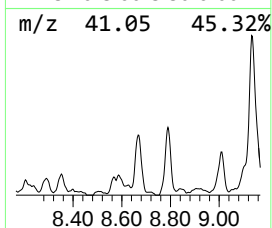
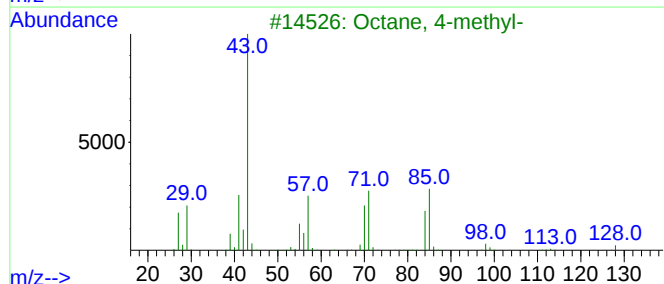
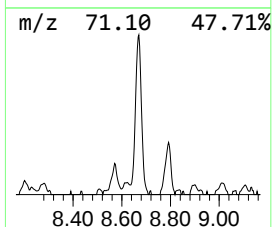
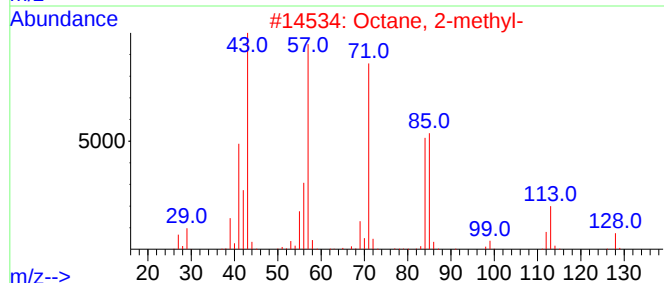
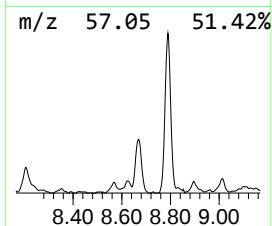
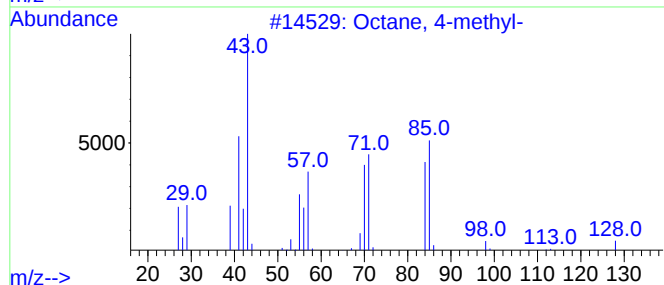
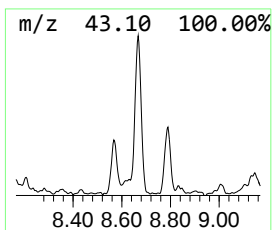
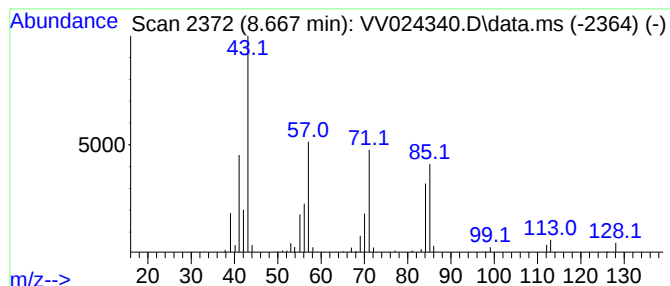
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.667	6.44 ug/L	92449	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-methyl-			128	C9H20	002216-34-4	90
2	Octane, 2-methyl-			128	C9H20	003221-61-2	81
3	Octane, 4-methyl-			128	C9H20	002216-34-4	74
4	Octane, 2-methyl-			128	C9H20	003221-61-2	74
5	Octane, 2-methyl-			128	C9H20	003221-61-2	59



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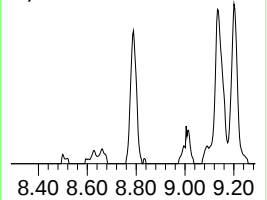
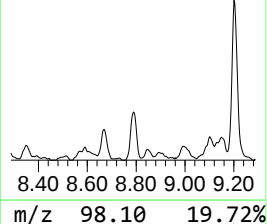
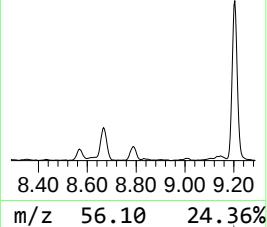
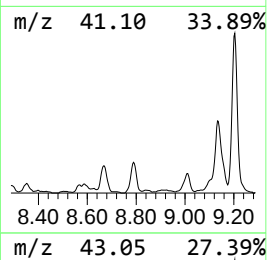
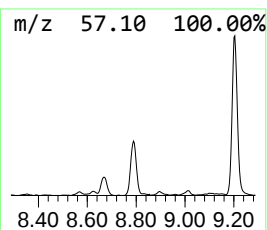
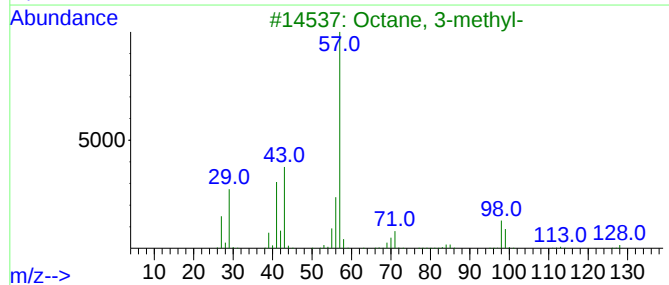
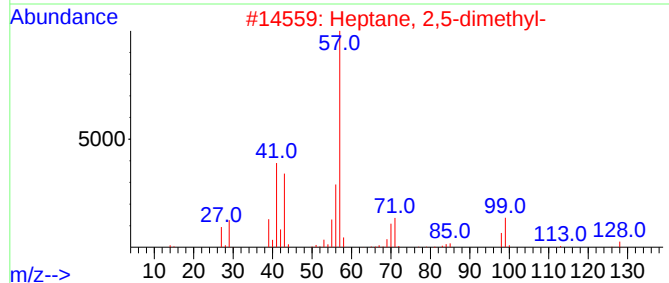
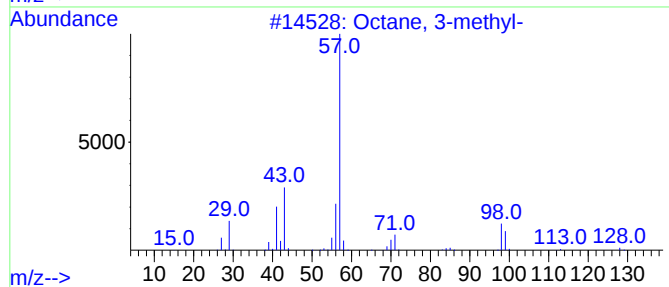
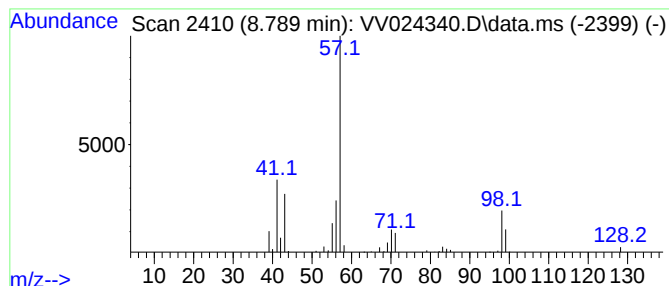
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Quant Title : VOC Analysis

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.789	6.12 ug/L	87925	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-			128	C9H20	002216-33-3	86
2	Heptane, 2,5-dimethyl-			128	C9H20	002216-30-0	81
3	Octane, 3-methyl-			128	C9H20	002216-33-3	72
4	Undecane, 6-methyl-			170	C12H26	017302-33-9	53
5	Octane, 3-methyl-			128	C9H20	002216-33-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

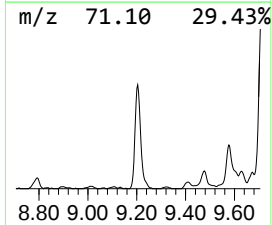
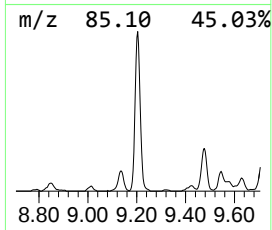
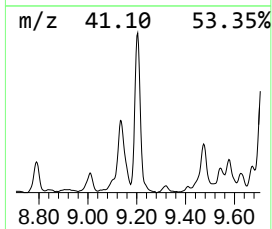
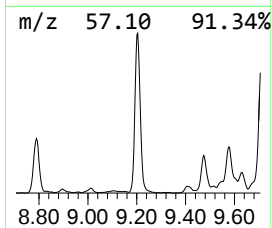
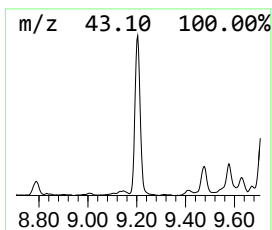
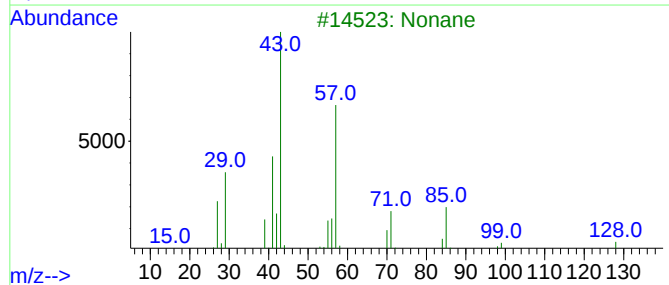
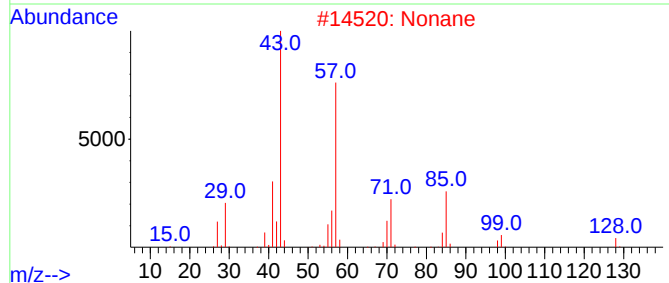
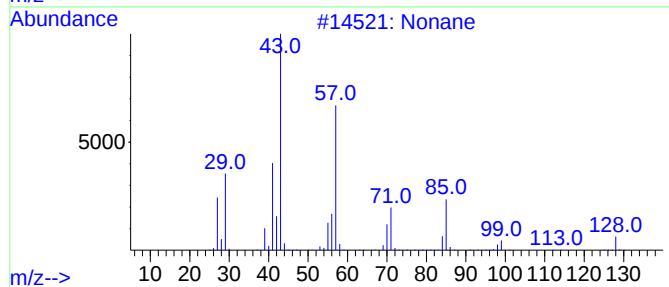
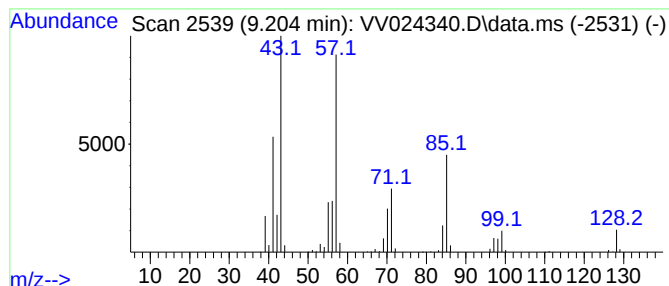
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Quant Title : VOC Analysis

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.204	34.98 ug/L	502441	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	94
2	Nonane			128	C9H20	000111-84-2	94
3	Nonane			128	C9H20	000111-84-2	94
4	Nonane			128	C9H20	000111-84-2	90
5	Octane			114	C8H18	000111-65-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

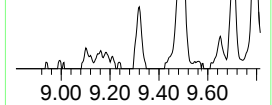
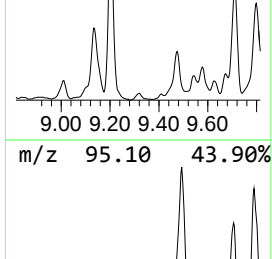
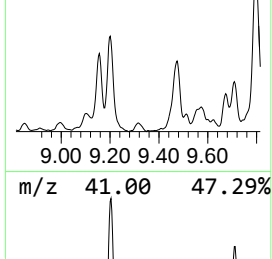
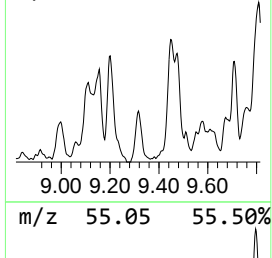
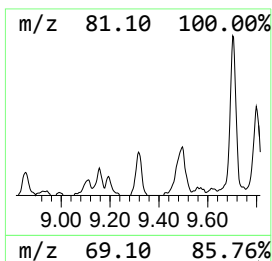
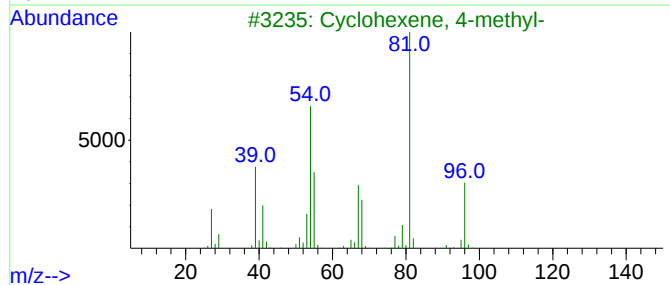
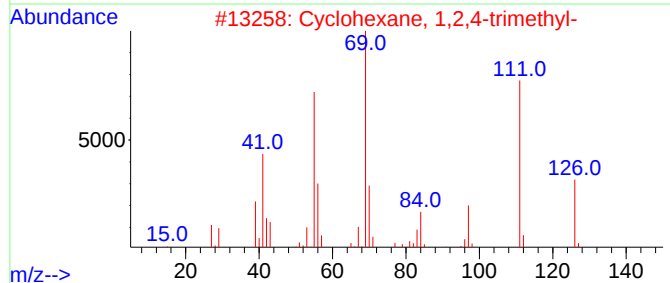
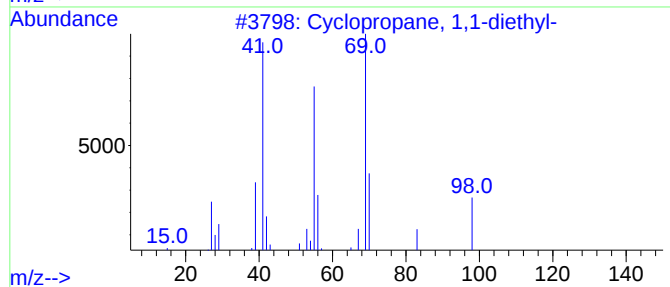
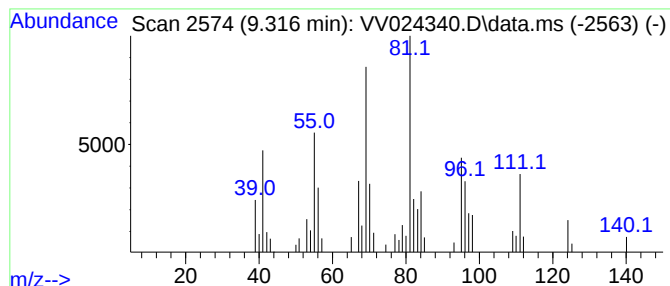
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Quant Title : VOC Analysis

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

Peak Number 6 (DEL) Alkane: Cyclic9.316 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.316	2.73 ug/L	39203	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,1-diethyl-	98	C7H14	001003-19-6	35
2			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	27
3			Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	25
4			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	25
5			Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

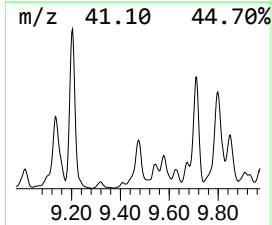
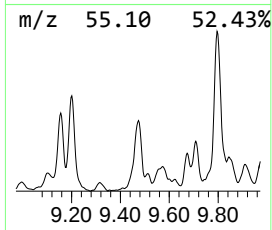
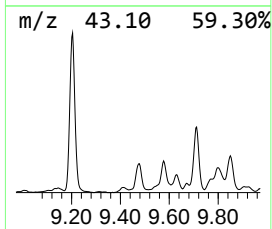
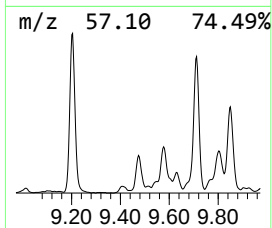
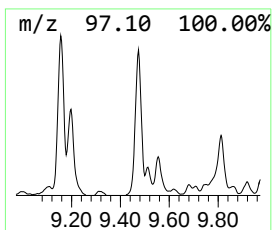
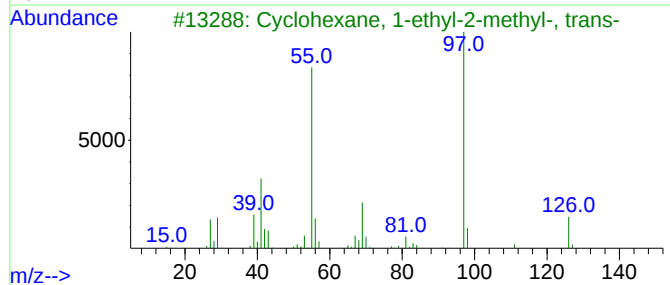
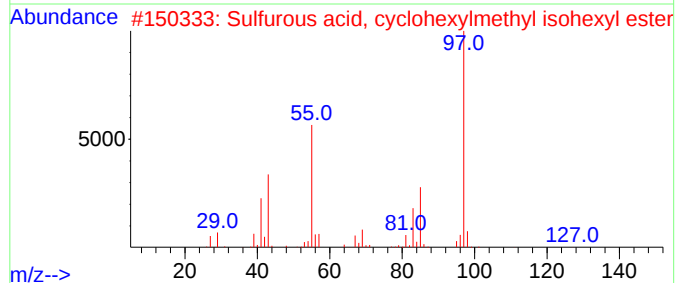
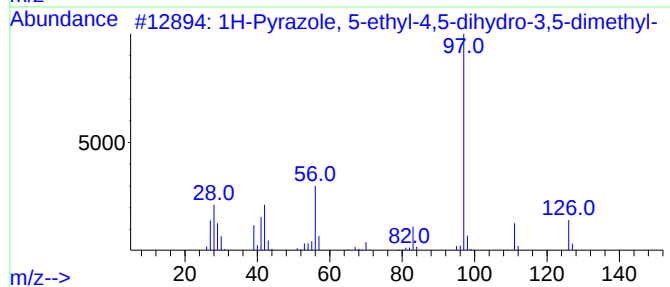
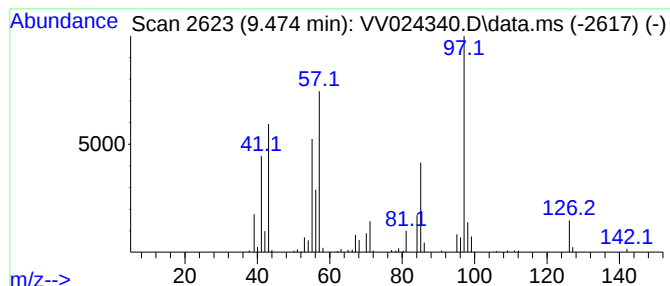
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Quant Title : VOC Analysis

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

Peak Number 7 unknown-01 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.474	9.69 ug/L	139252	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrazole, 5-ethyl-4,5-dihydro...	126	C7H14N2	021981-22-6	49
2			Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	47
3			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	46
4			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	43
5			4-Octen-3-one	126	C8H14O	014129-48-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

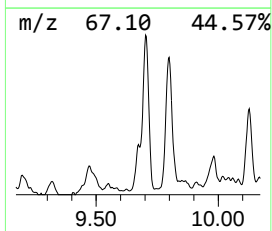
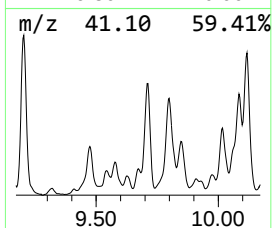
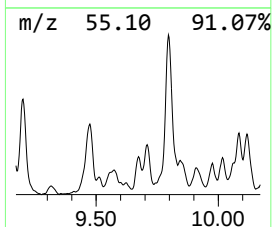
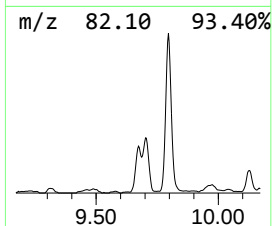
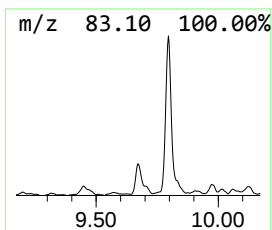
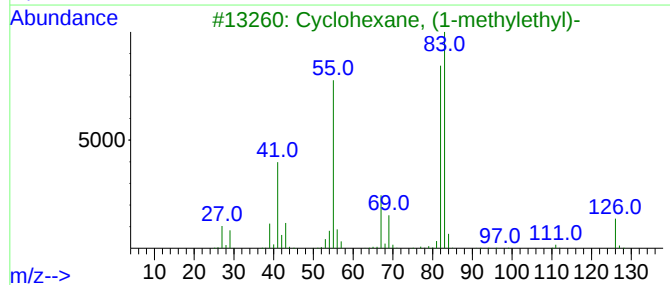
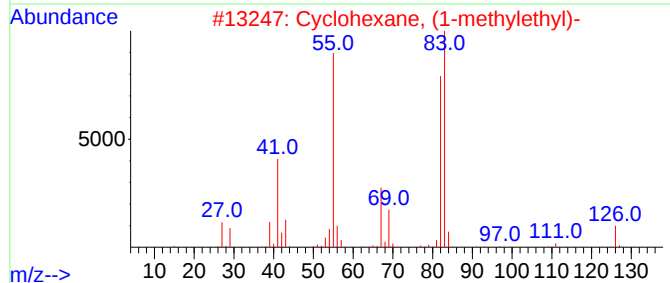
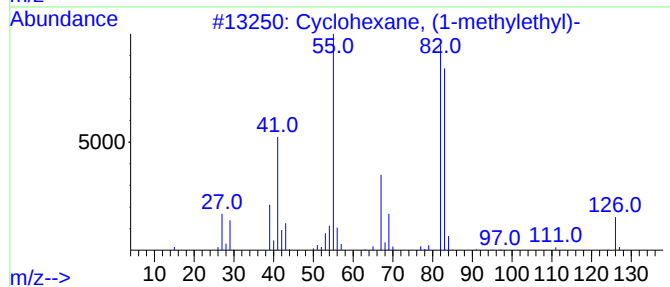
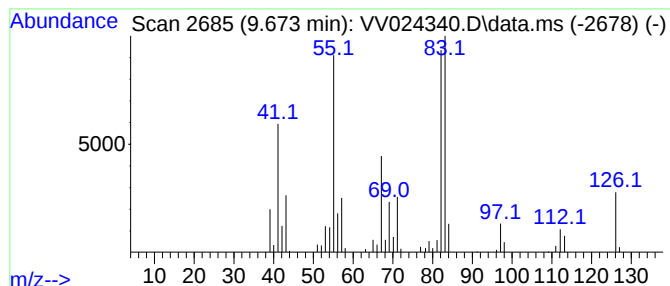
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Quant Title : VOC Analysis

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

Peak Number 8 (DEL) Alkane: Cyclic9.673 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.673	3.91 ug/L	56093	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	87
2			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	76
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	76
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	68
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

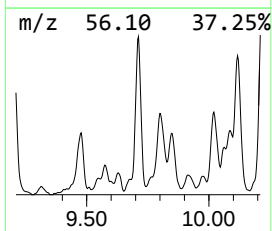
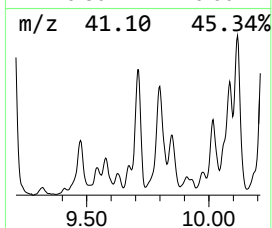
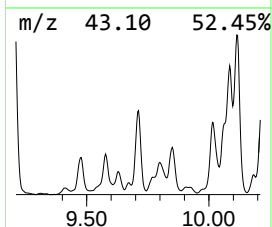
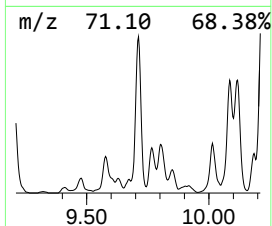
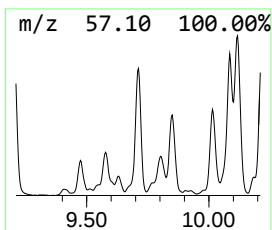
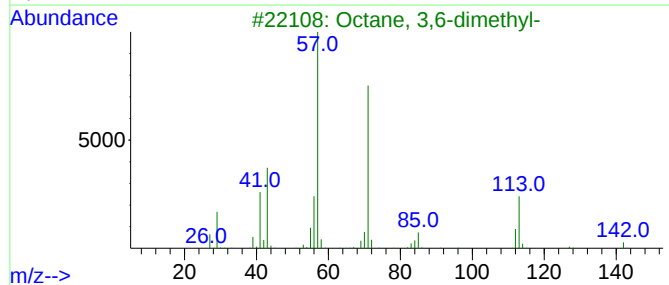
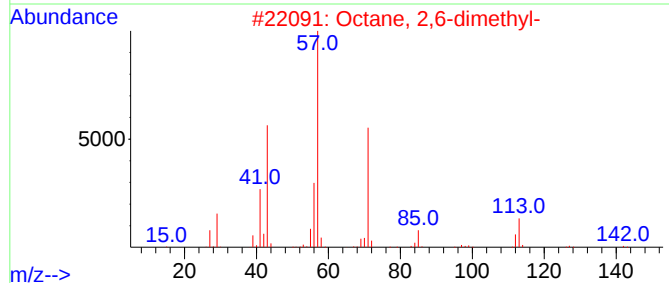
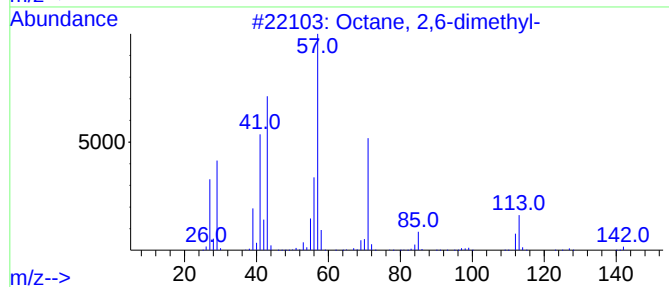
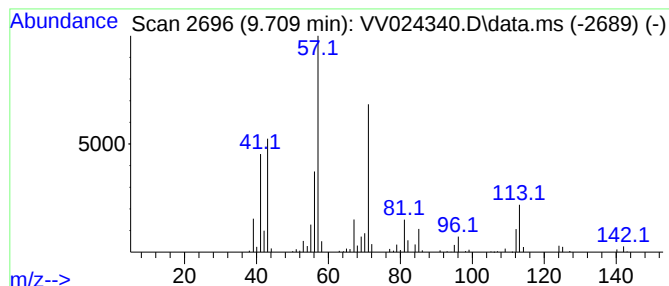
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Quant Title : VOC Analysis

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.709	24.11 ug/L	346295	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	90
2	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	90
3	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	90
4	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	83
5	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

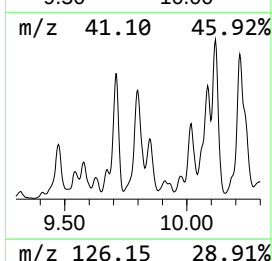
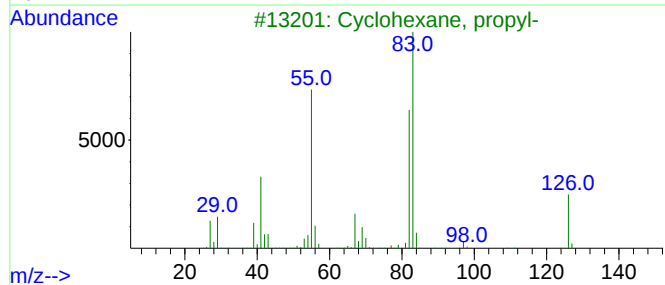
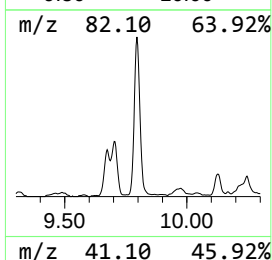
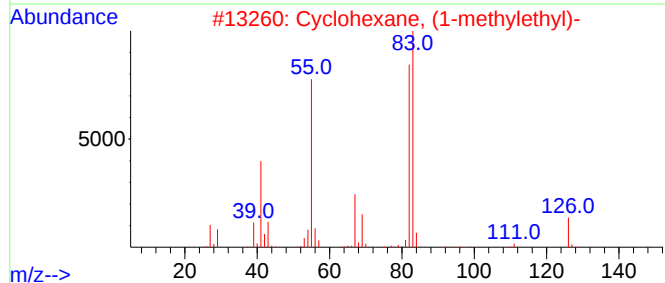
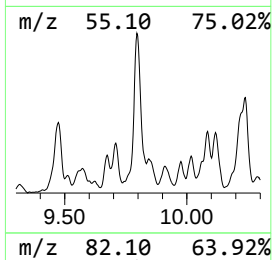
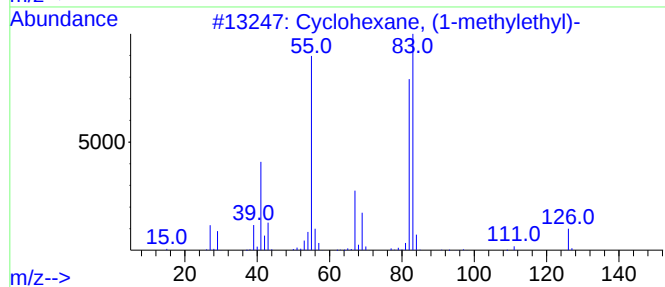
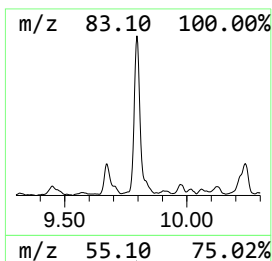
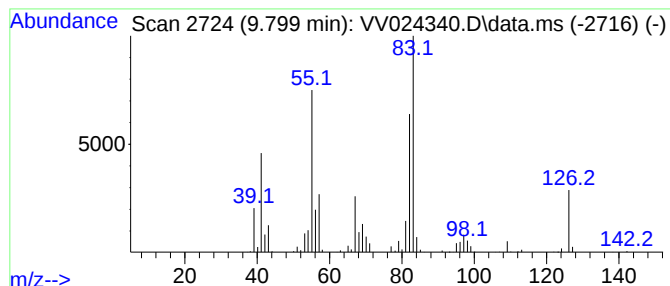
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Quant Title : VOC Analysis

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

Peak Number 10 (DEL) Alkane: Cyclic9.799 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.799	20.11 ug/L	288880	Chlorobenzene-d5	8.850

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	87
2		Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	87
3		Cyclohexane, propyl-	126	C9H18	001678-92-8	81
4		Cyclohexane, propyl-	126	C9H18	001678-92-8	64
5		Cyclohexane, propyl-	126	C9H18	001678-92-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

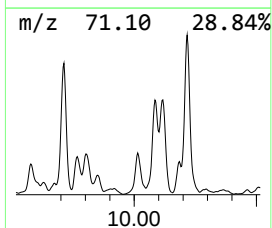
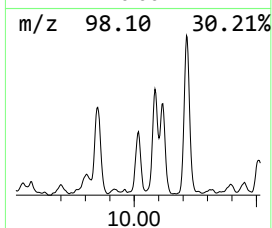
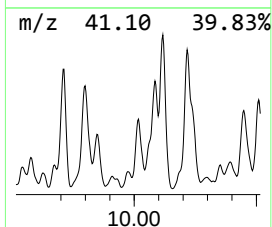
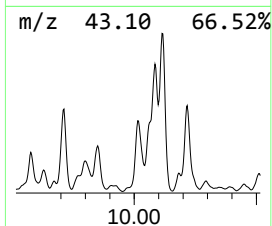
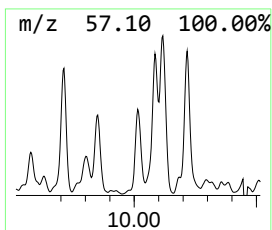
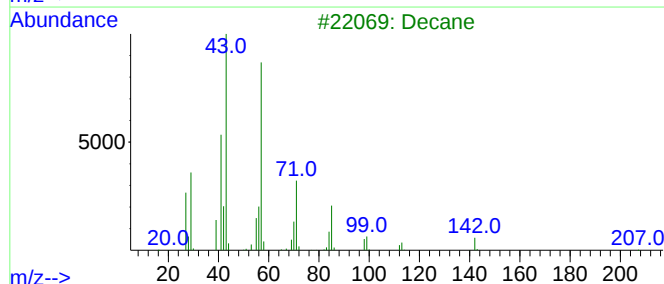
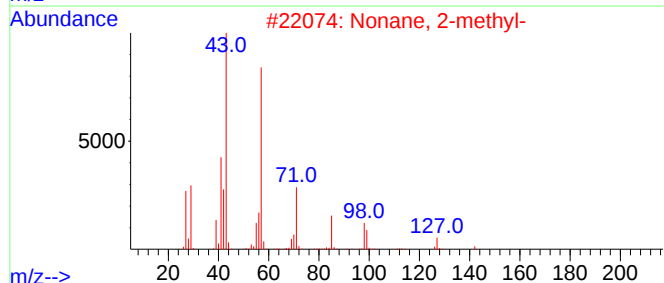
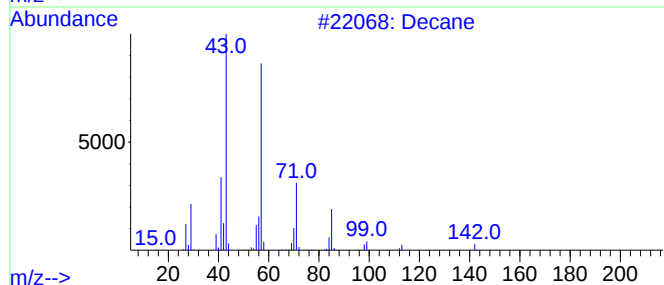
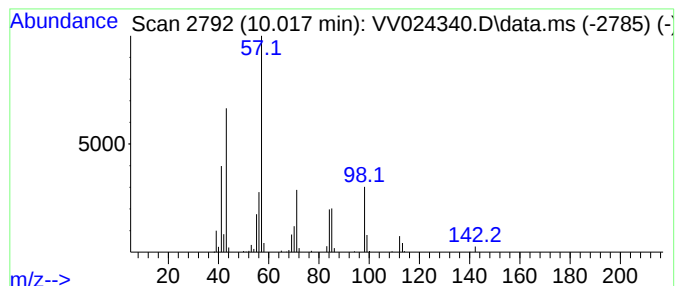
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Quant Title : VOC Analysis

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.017	9.82 ug/L	140992	Chlorobenzene-d5	8.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	52
2			Nonane, 2-methyl-	142	C10H22	000871-83-0	50
3			Decane	142	C10H22	000124-18-5	47
4			Octane, 2,2,6-trimethyl-	156	C11H24	062016-28-8	47
5			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

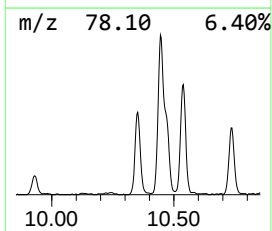
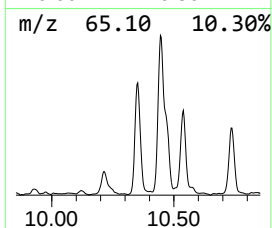
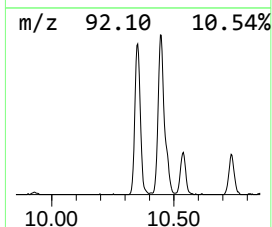
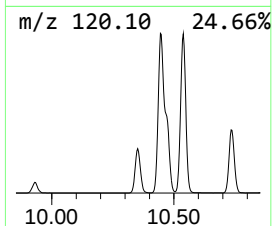
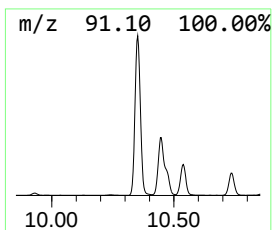
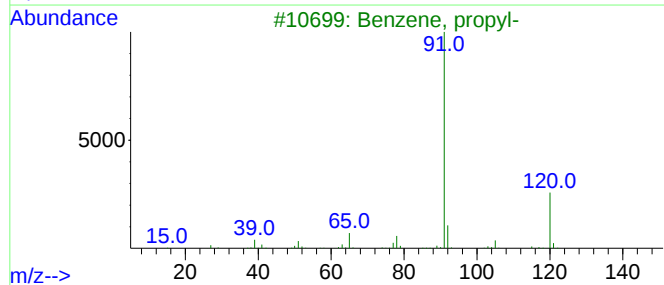
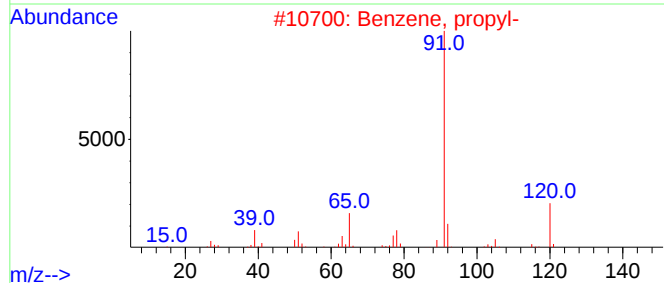
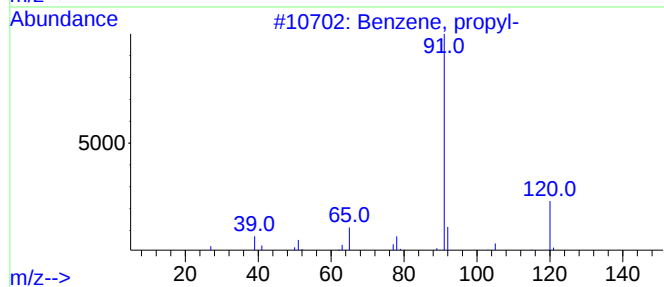
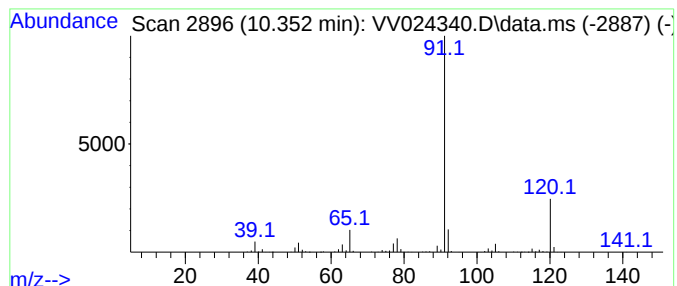
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.352	32.60 ug/L	607309	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	93
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			Benzene, propyl-	120	C9H12	000103-65-1	90
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

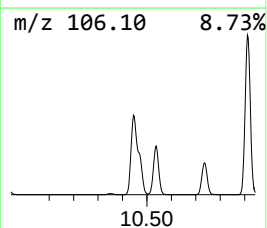
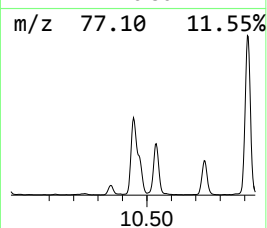
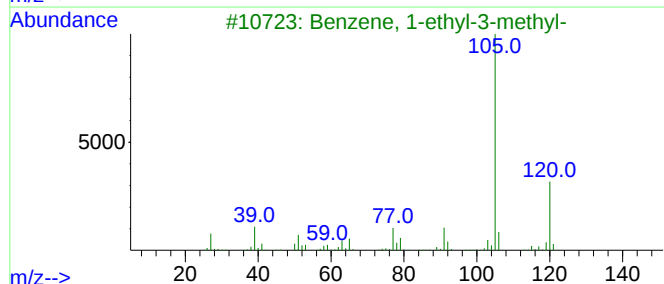
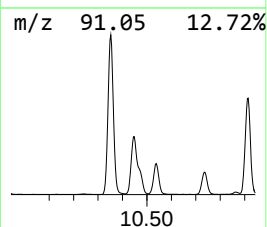
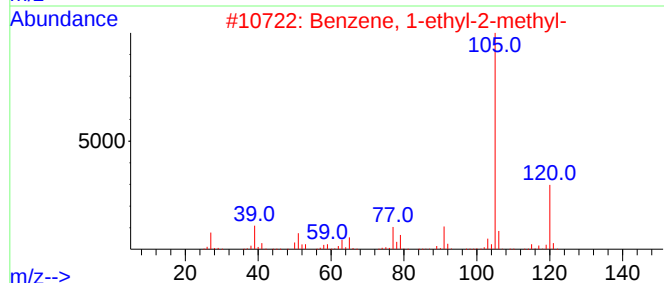
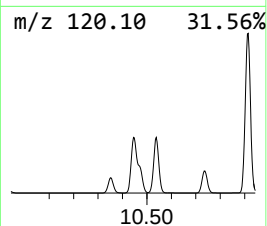
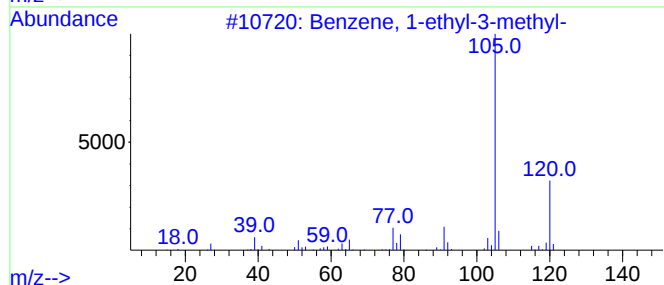
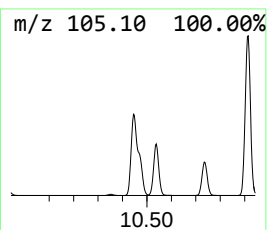
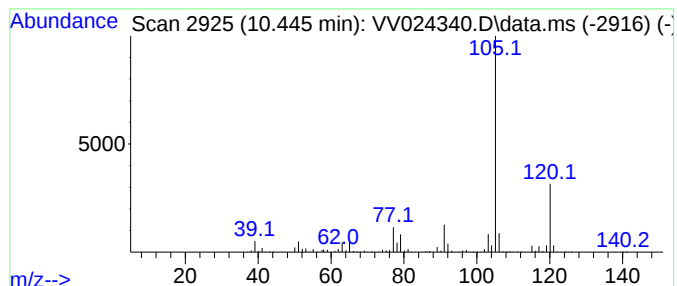
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Quant Title : VOC Analysis

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

Peak Number 13 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.445	161.05 ug/L	3000180	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

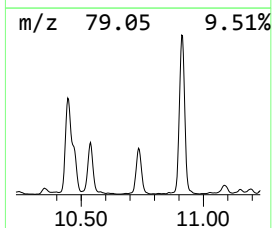
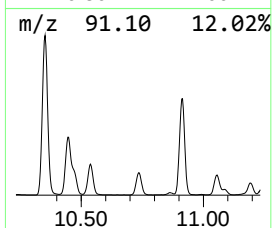
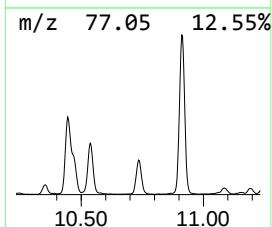
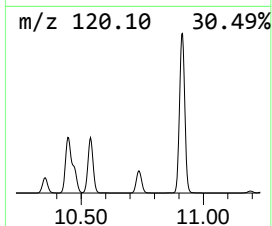
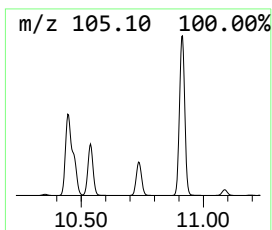
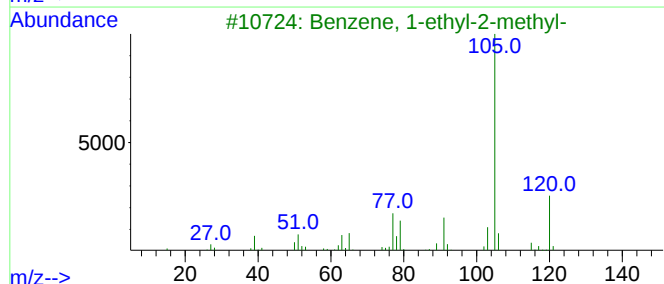
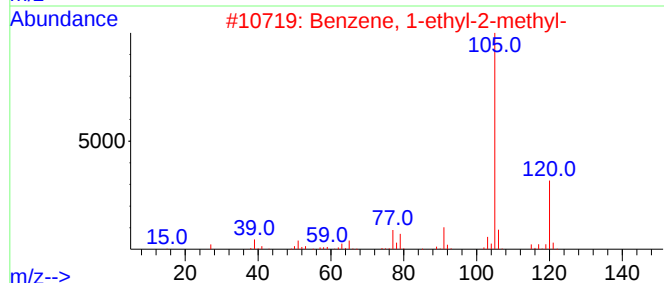
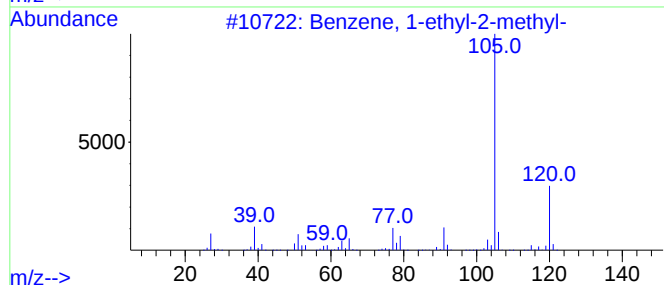
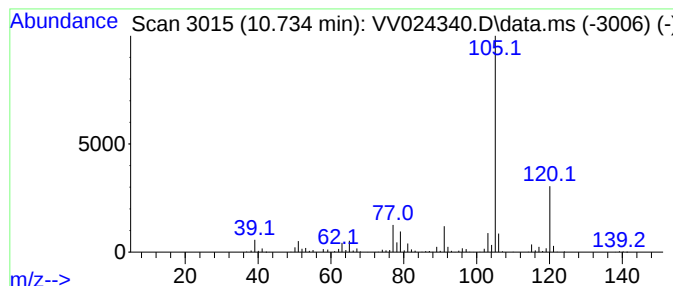
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Quant Title : VOC Analysis

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

Peak Number 14 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.734	57.22 ug/L	1065950	1,4-Dichlorobenzene-d4	11.249

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-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

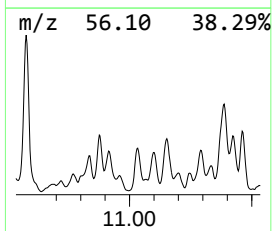
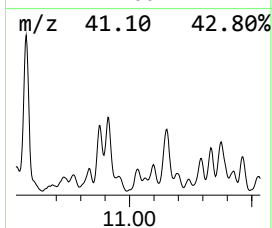
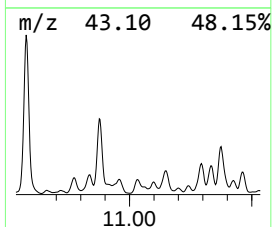
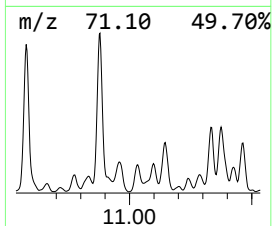
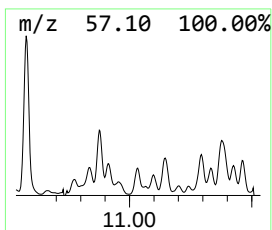
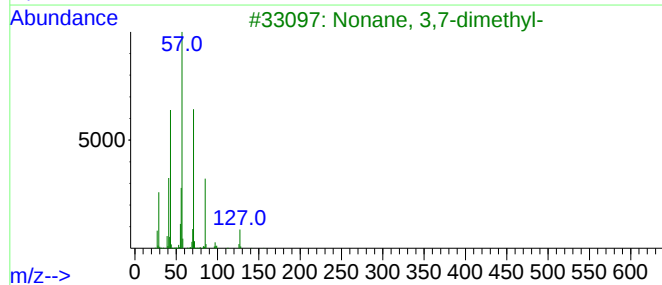
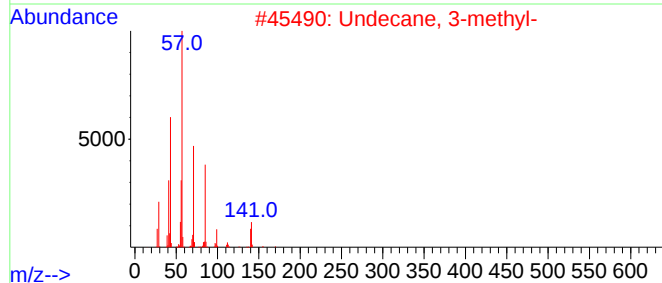
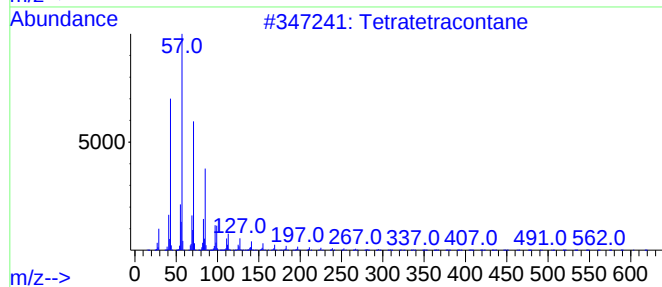
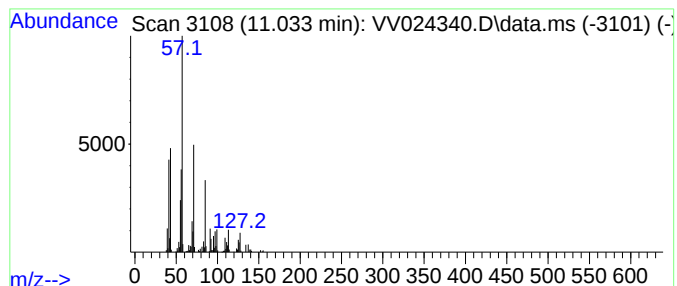
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Quant Title : VOC Analysis

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.034	8.55 ug/L	159293	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	59
2			Undecane, 3-methyl-	170	C12H26	001002-43-3	59
3			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	50
4			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	50
5			Dodecane, 3-methyl-	184	C13H28	017312-57-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

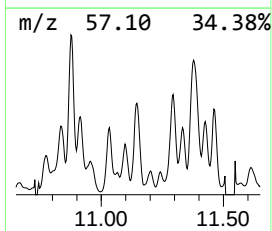
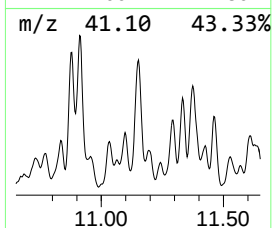
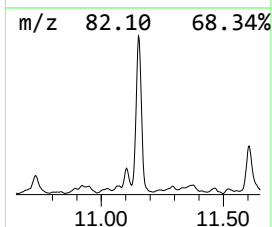
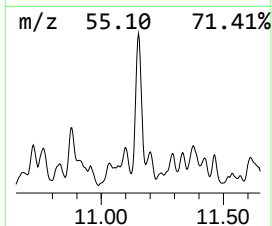
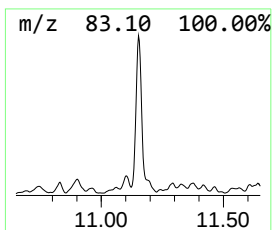
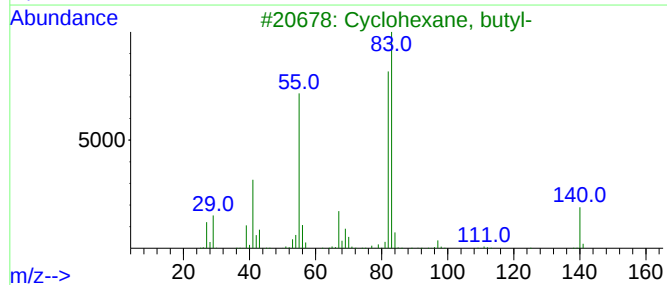
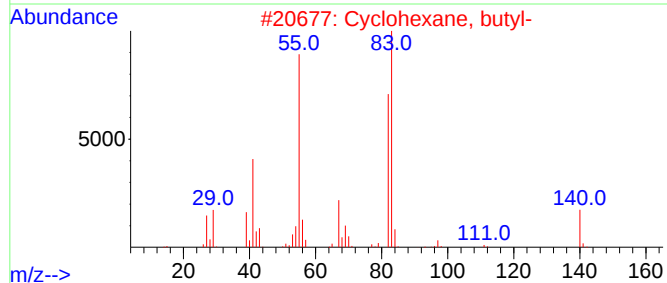
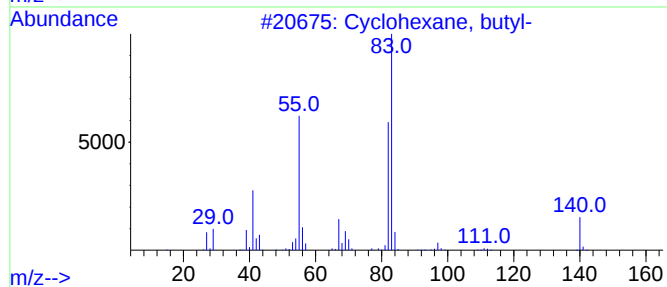
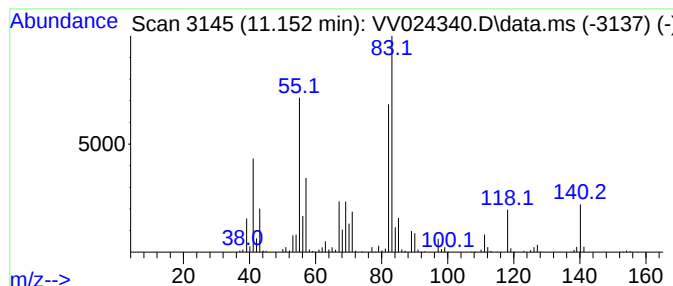
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Quant Title : VOC Analysis

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

Peak Number 16 (DEL) Alkane: Cyclic11.152 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.152	34.38 ug/L	640398	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	70
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	64
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	58
4			n-Heptadecylcyclohexane	322	C23H46	019781-73-8	53
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

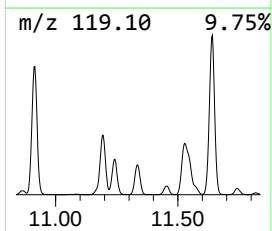
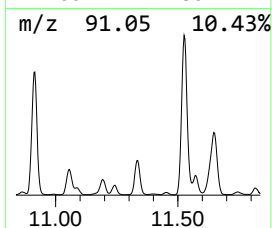
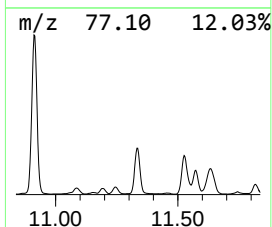
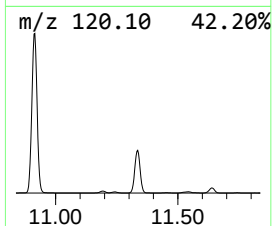
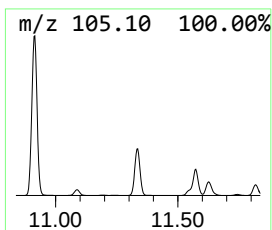
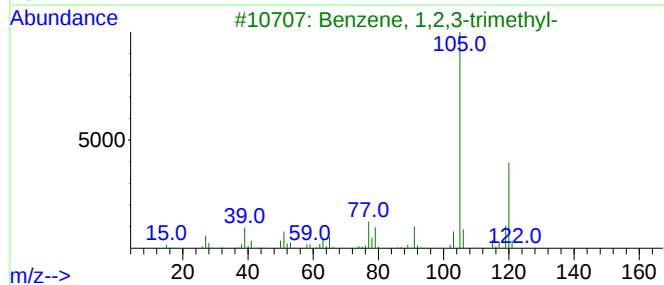
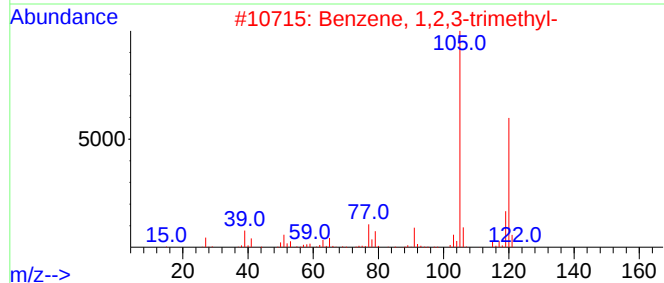
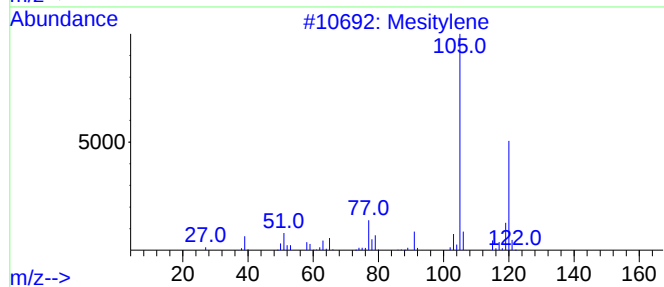
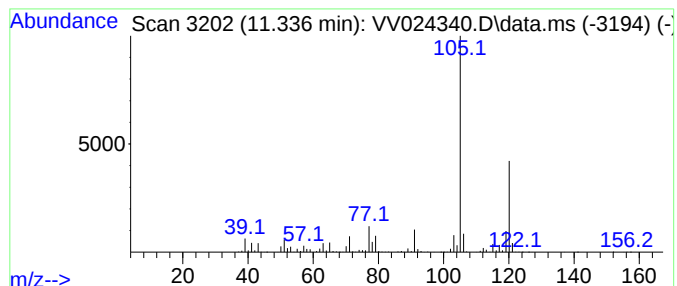
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 17 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	68.75 ug/L	1280660	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

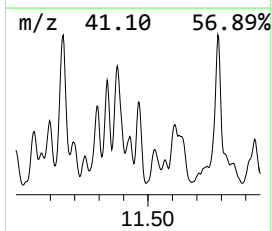
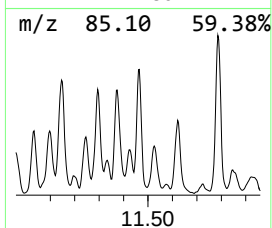
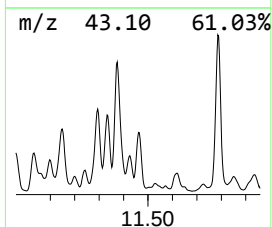
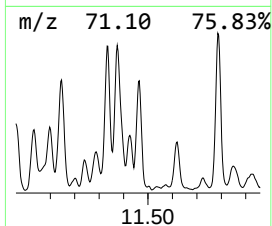
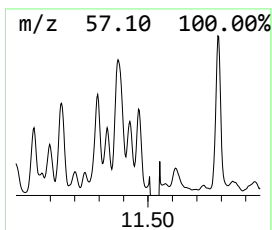
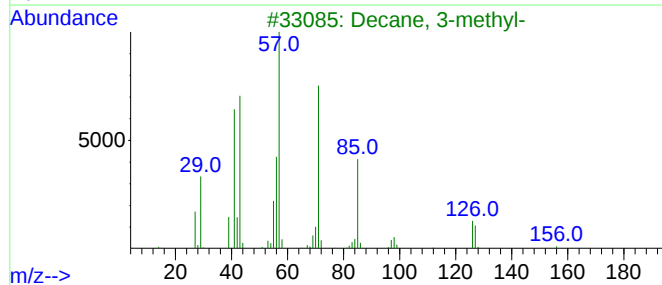
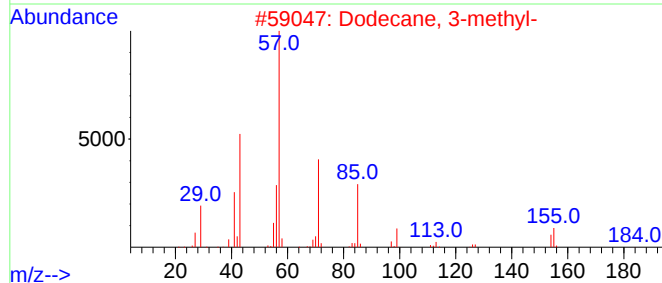
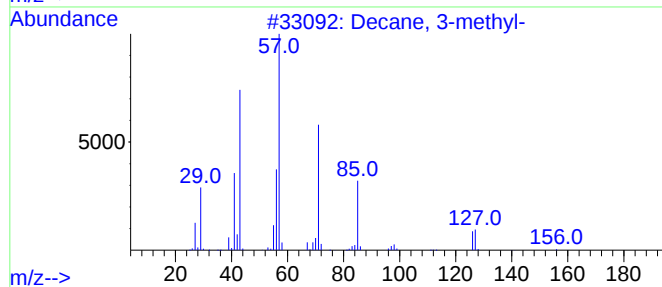
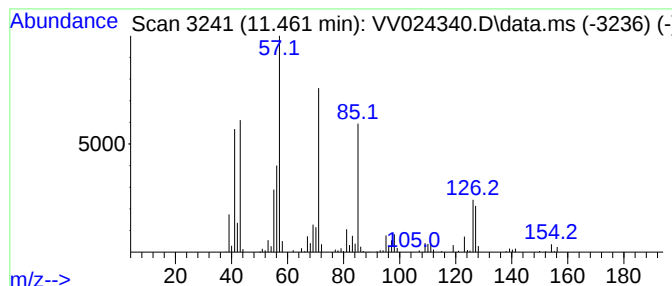
Instrument :
MSVOA_V
ClientSampleId :
BGLN2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.461	13.36 ug/L	248818	1,4-Dichlorobenzene-d4			11.249
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-		156	C11H24	013151-34-3	93
2	Dodecane, 3-methyl-		184	C13H28	017312-57-1	53
3	Decane, 3-methyl-		156	C11H24	013151-34-3	50
4	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	47
5	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

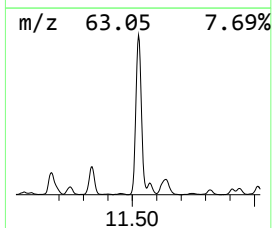
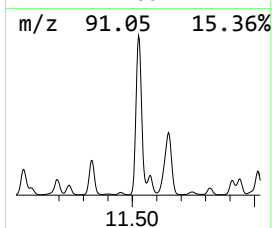
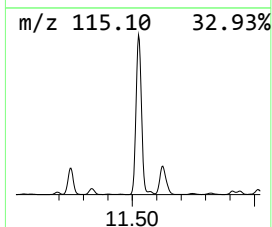
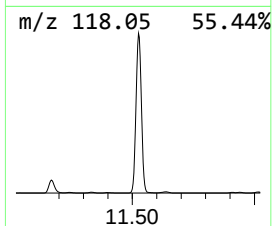
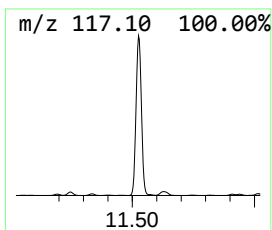
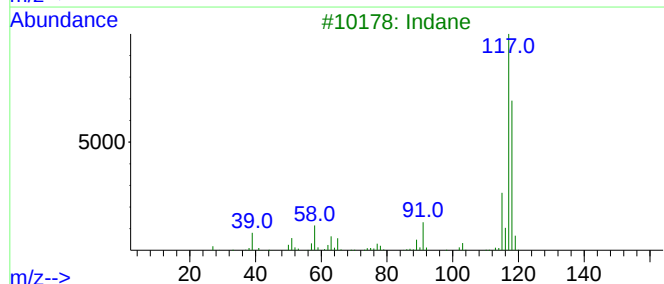
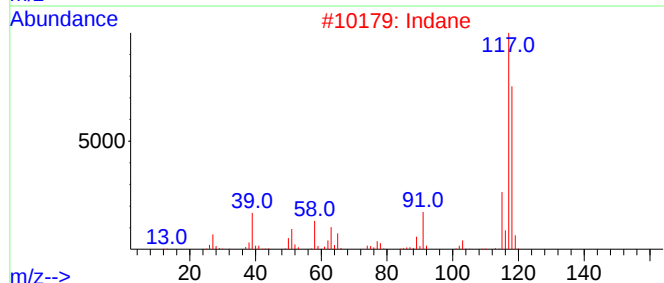
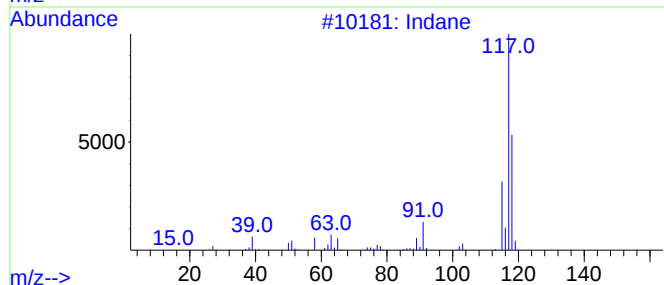
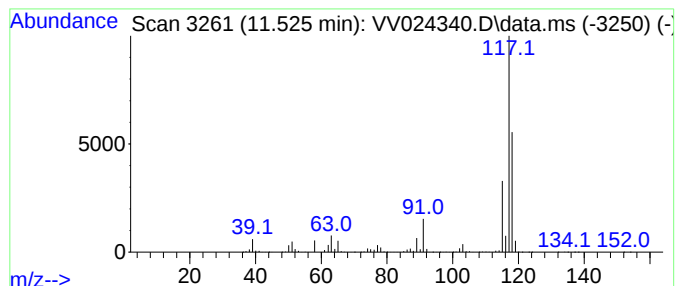
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Quant Title : VOC Analysis

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

Peak Number 19 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.525	253.84 ug/L	4728670	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Indane	118	C9H10	000496-11-7	91
3			Indane	118	C9H10	000496-11-7	87
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

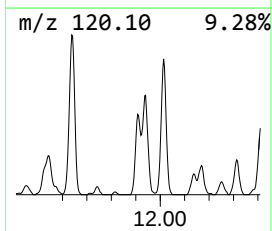
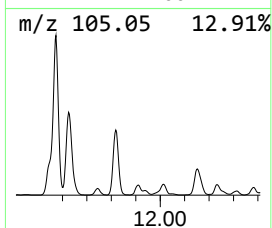
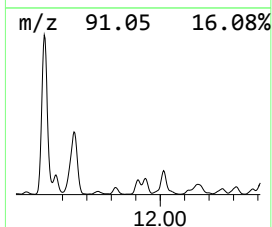
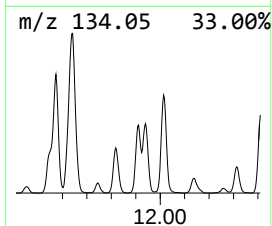
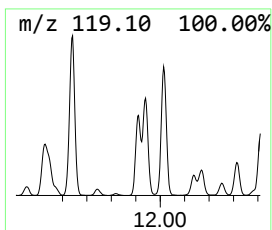
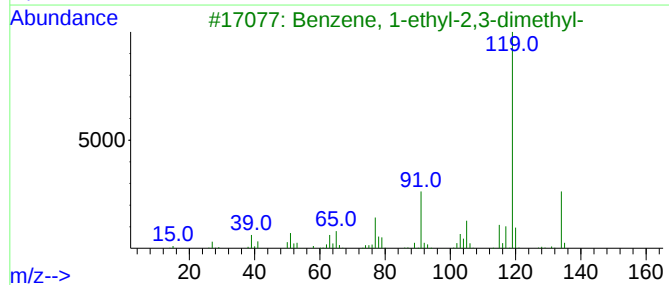
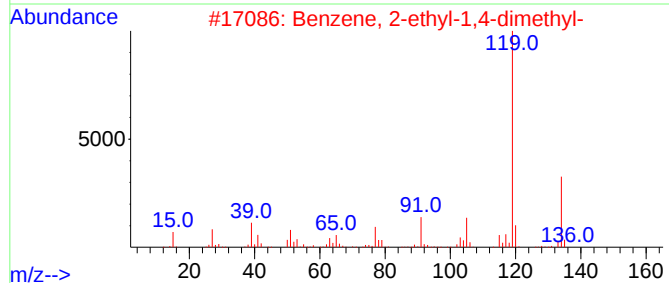
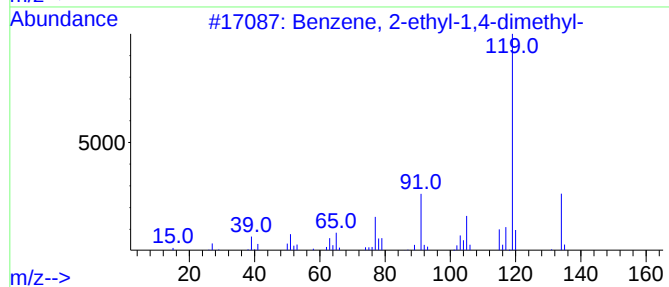
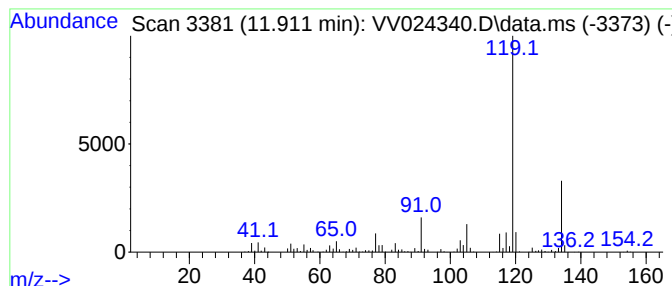
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Quant Title : VOC Analysis

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

Peak Number 20 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	17.57 ug/L	327390	1,4-Dichlorobenzene-d4	11.249

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

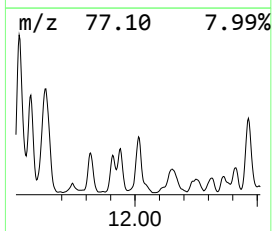
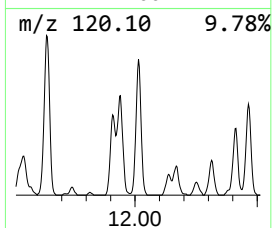
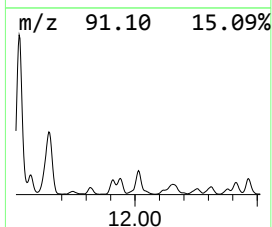
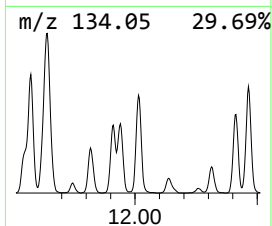
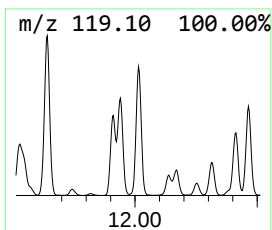
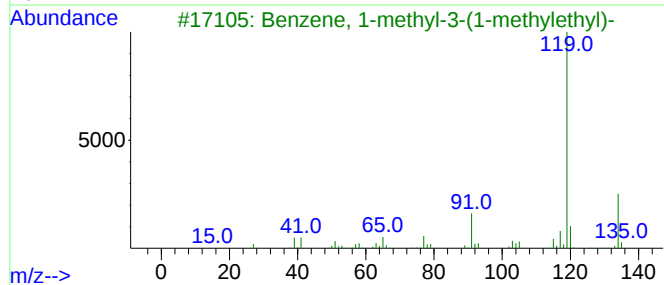
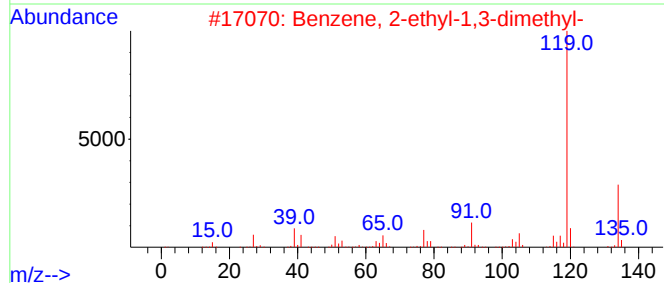
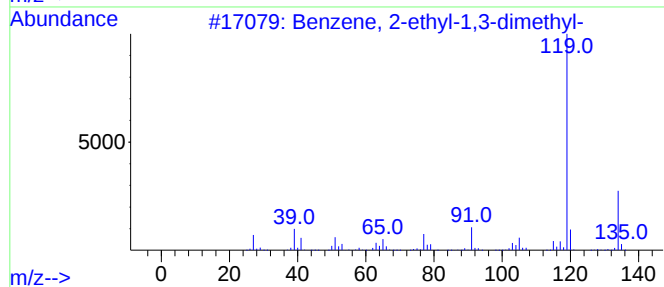
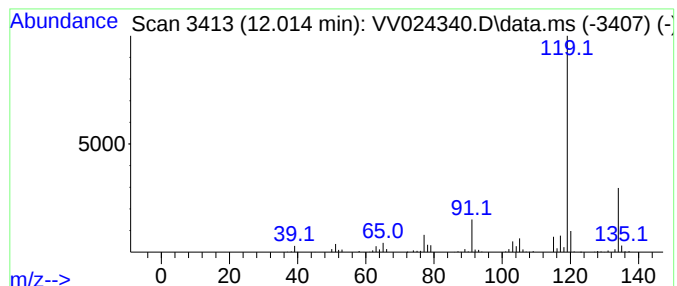
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Quant Title : VOC Analysis

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

Peak Number 21 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.014	34.33 ug/L	639559	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

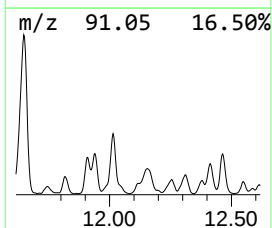
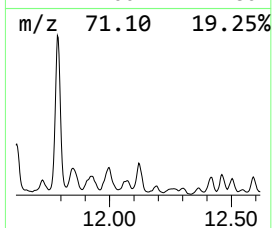
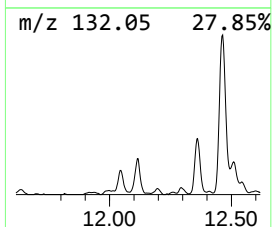
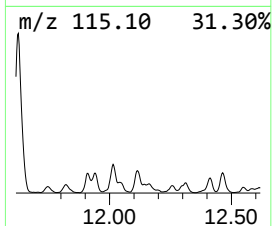
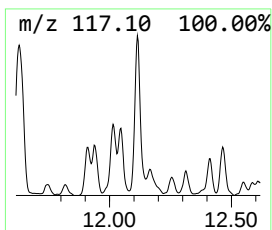
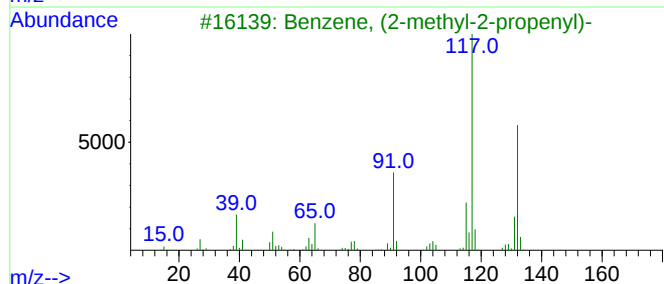
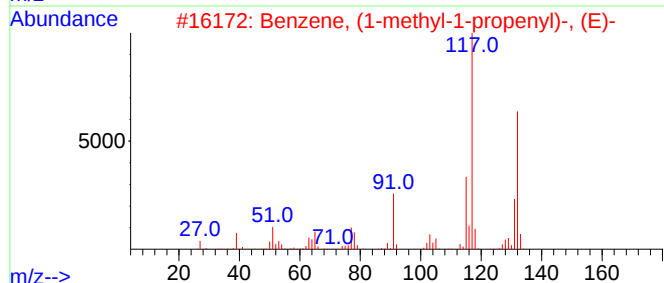
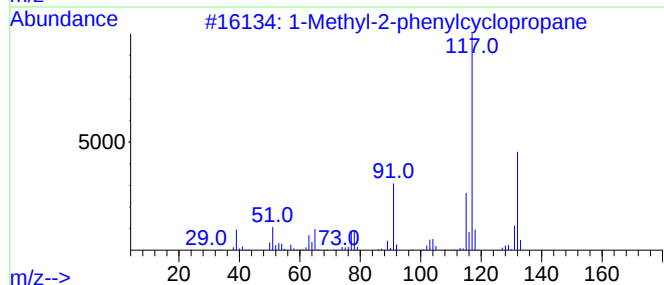
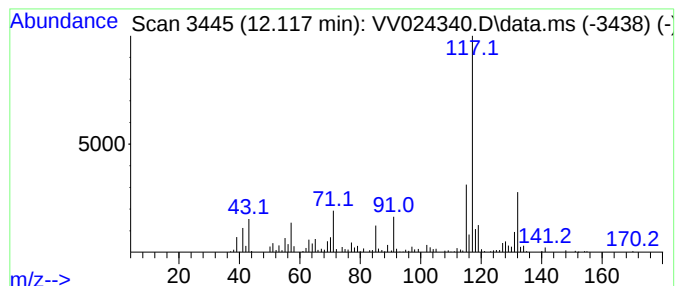
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Quant Title : VOC Analysis

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

Peak Number 22 1-Methyl-2-phenylcyclopropane Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.117	7.57 ug/L	140970	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	68
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	68
3			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	68
4			Indan, 1-methyl-	132	C10H12	000767-58-8	64
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

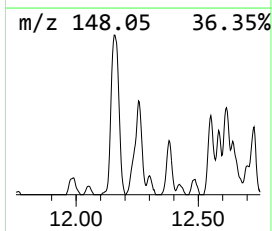
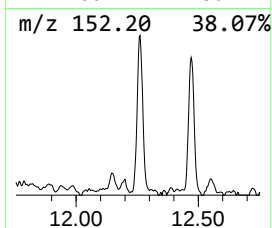
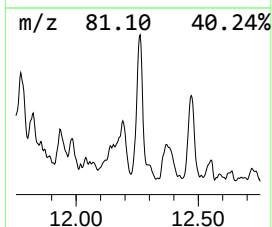
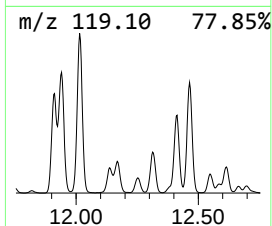
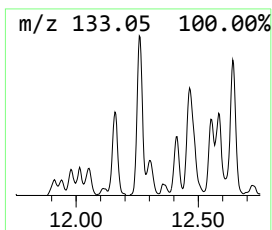
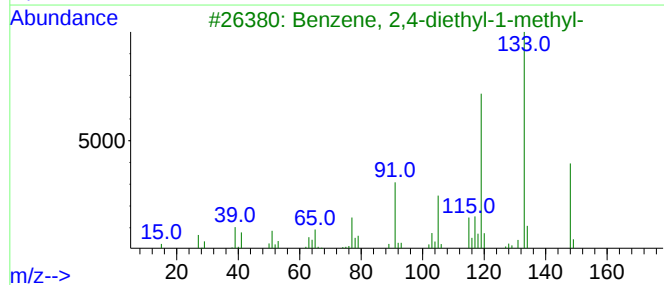
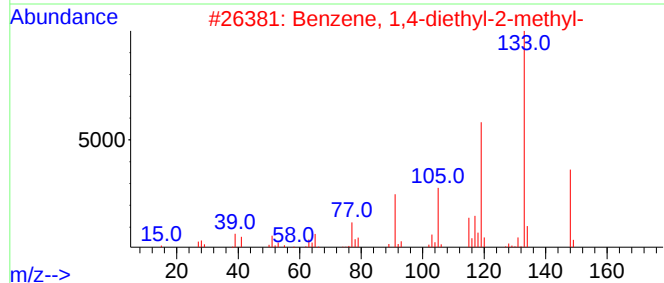
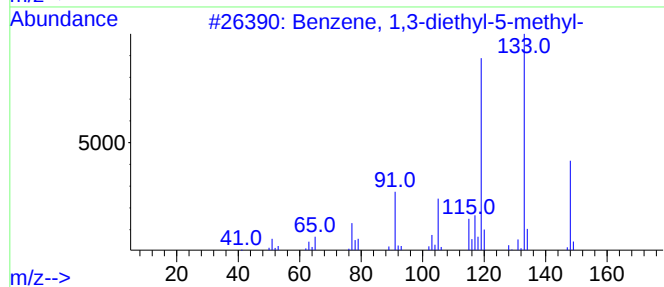
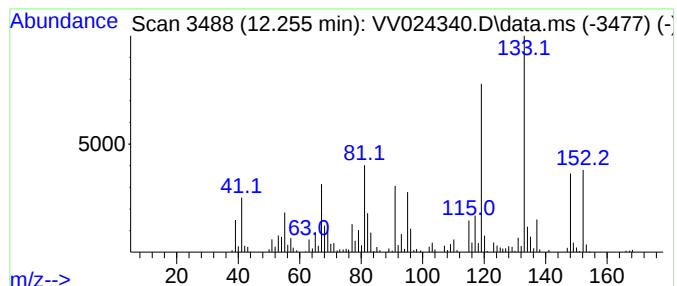
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Quant Title : VOC Analysis

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

Peak Number 23 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.255	13.24 ug/L	246701	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	86
2			Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	83
3			Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	64
4			Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	64
5			Benzene, pentamethyl-	148	C11H16	000700-12-9	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

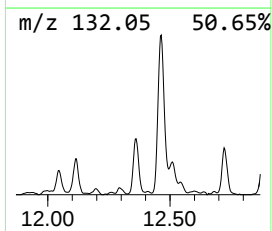
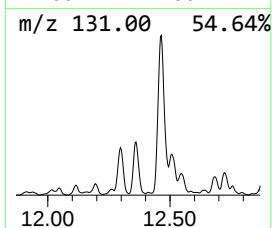
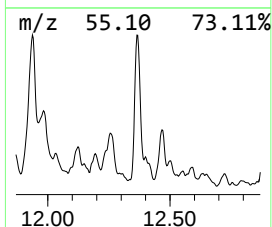
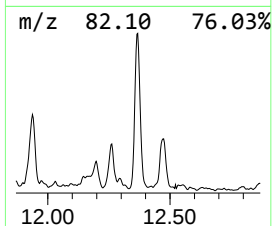
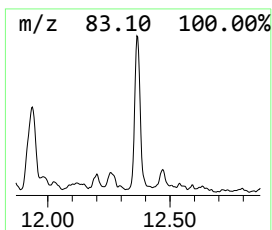
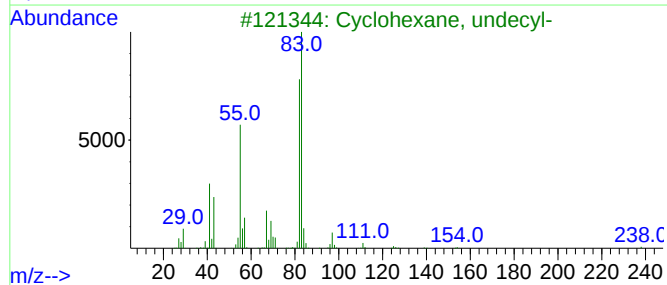
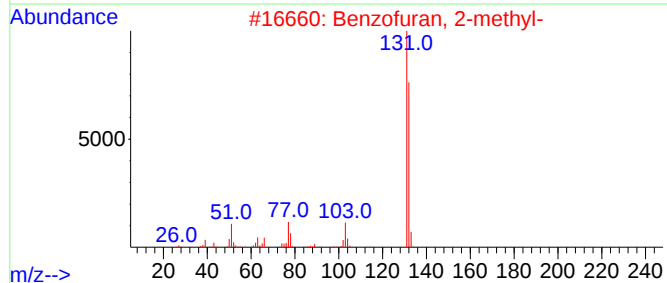
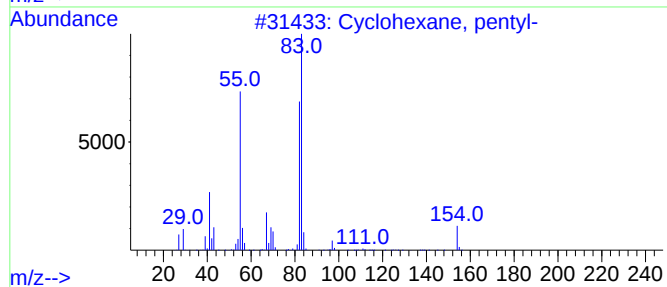
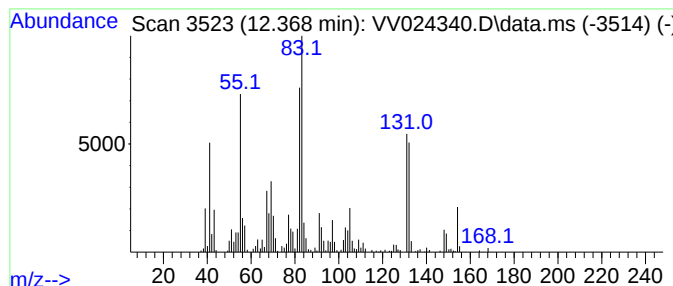
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Quant Title : VOC Analysis

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

Peak Number 24 (DEL) Alkane: Cyclic12.368 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.368	15.28 ug/L	284728	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	46
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	43
3			Cyclohexane, undecyl-	238	C17H34	054105-66-7	38
4			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	35
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

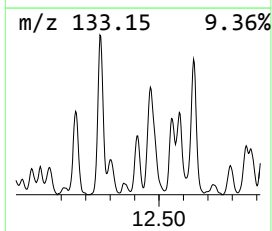
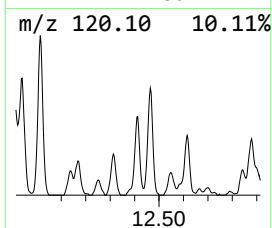
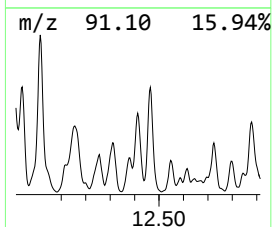
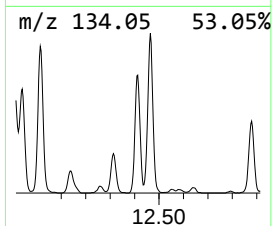
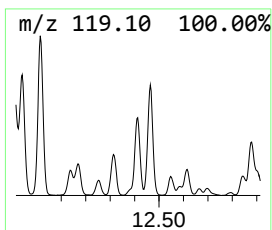
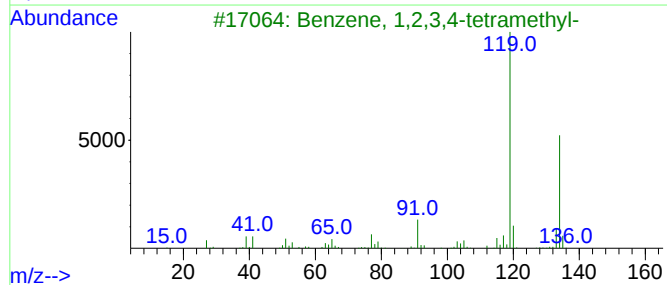
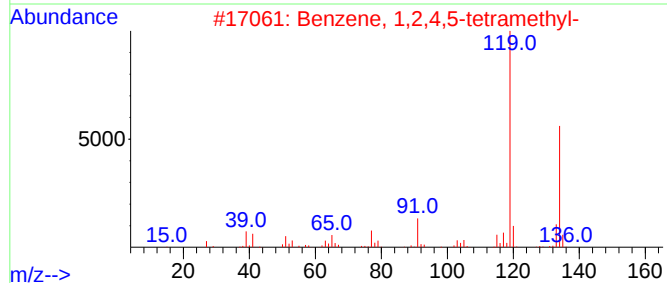
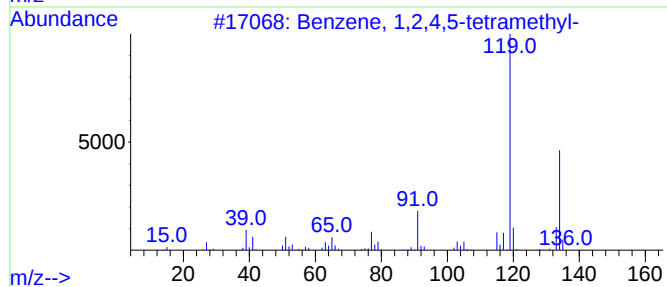
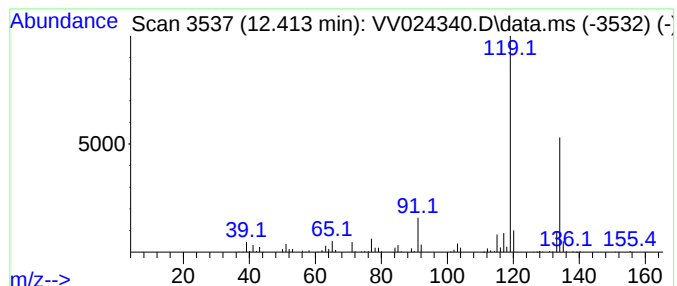
Instrument :
MSVOA_V
ClientSampleId :
BGLN2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
Quant Title : VOC Analysis

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

Peak Number 25 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.413	13.53 ug/L	252087	1,4-Dichlorobenzene-d4			11.249
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	95
2	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	94
3	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	94
4	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	94
5	Benzene, 1,2,3,5-tetramethyl-		134	C10H14	000527-53-7	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

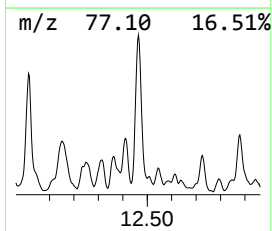
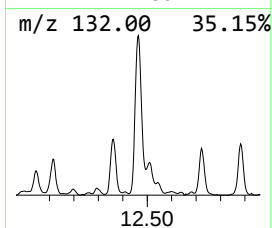
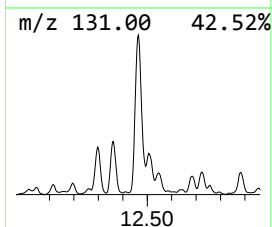
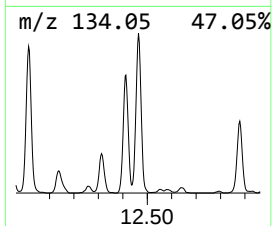
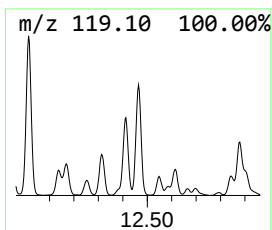
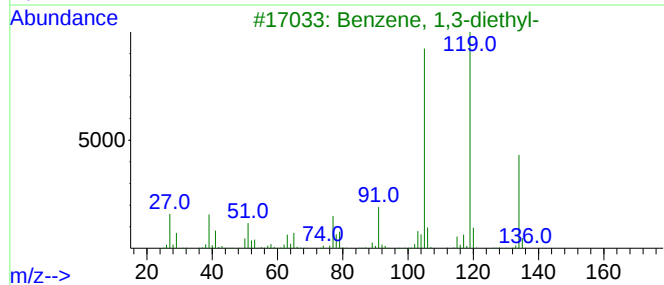
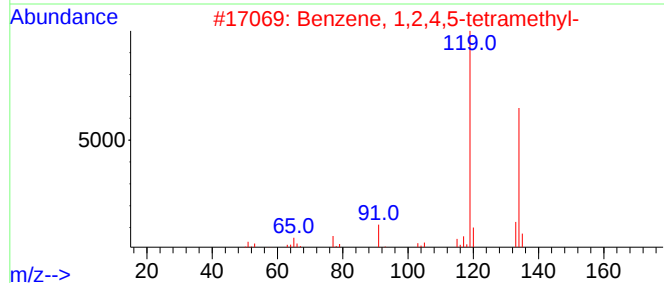
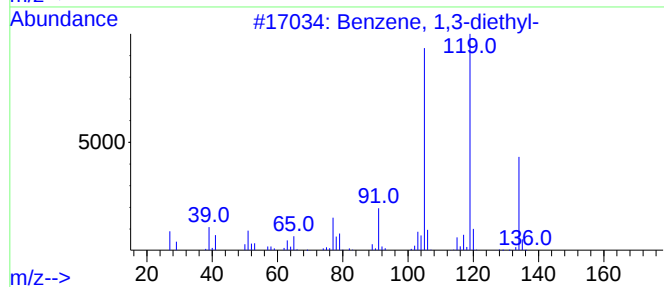
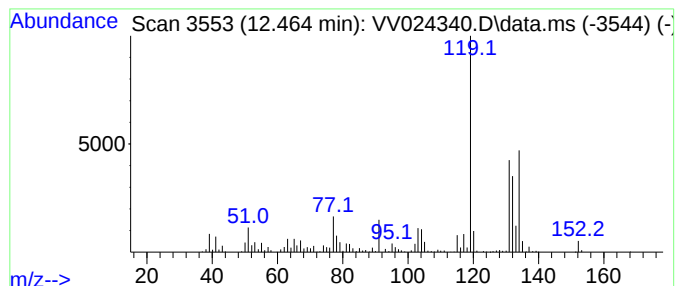
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Quant Title : VOC Analysis

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

Peak Number 26 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	36.97 ug/L	688778	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	64
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

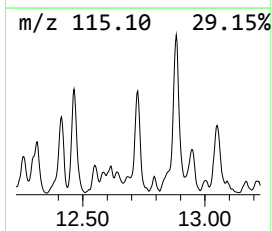
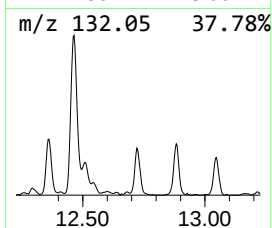
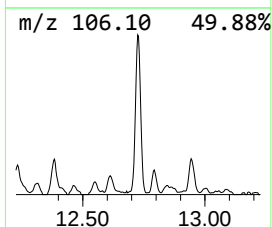
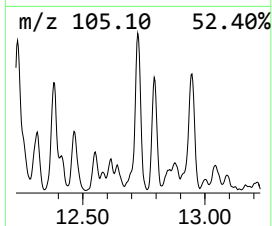
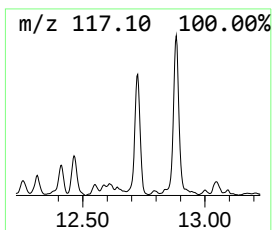
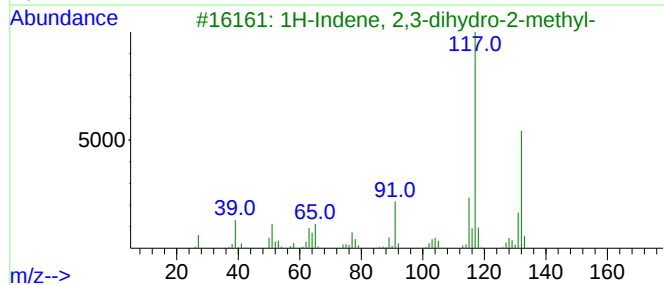
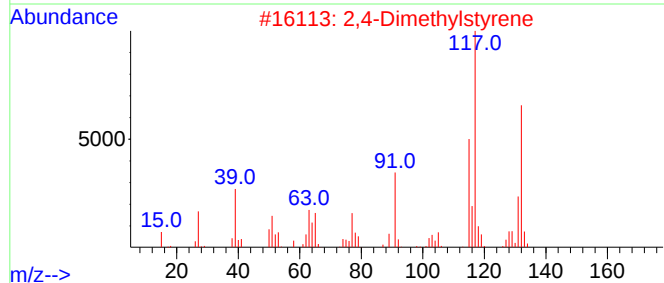
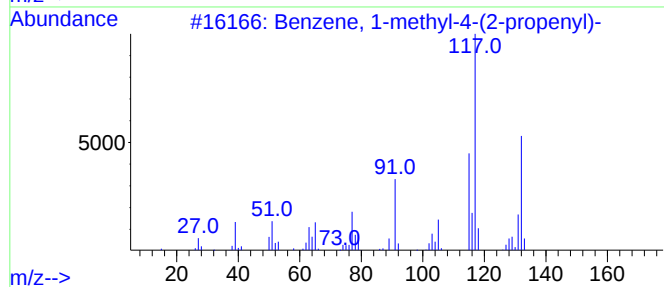
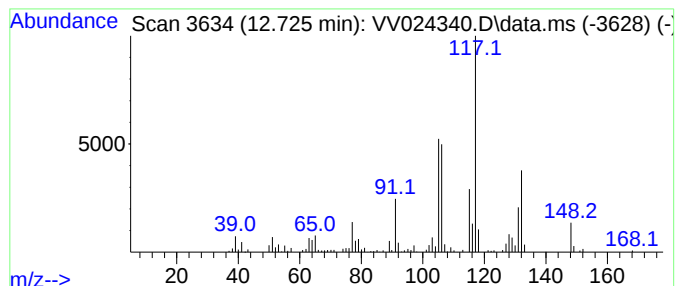
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Quant Title : VOC Analysis

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

Peak Number 27 Benzene, 1-methyl-4-(2-prop... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.725	10.46 ug/L	194836	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	87
3			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	70
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	70
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

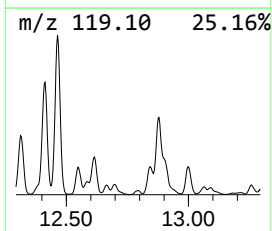
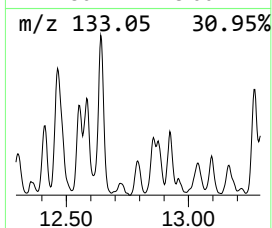
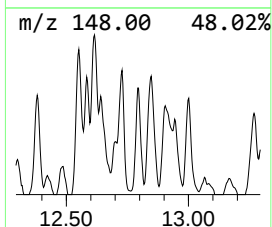
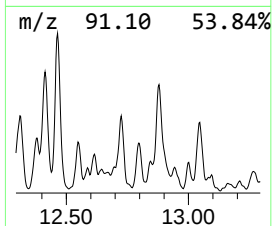
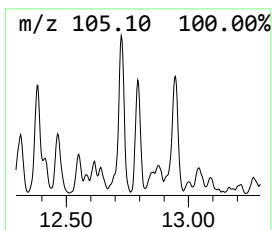
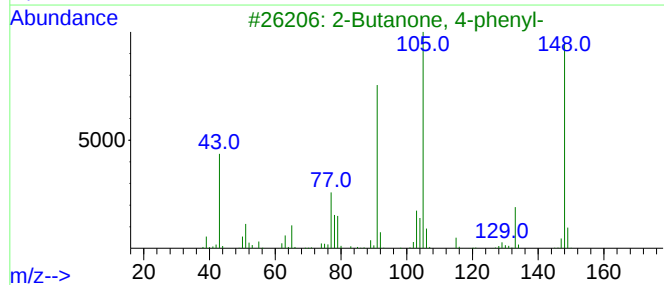
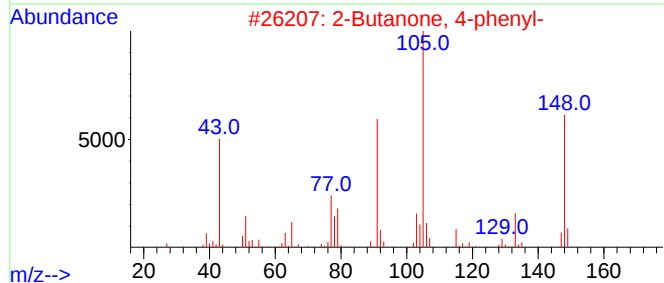
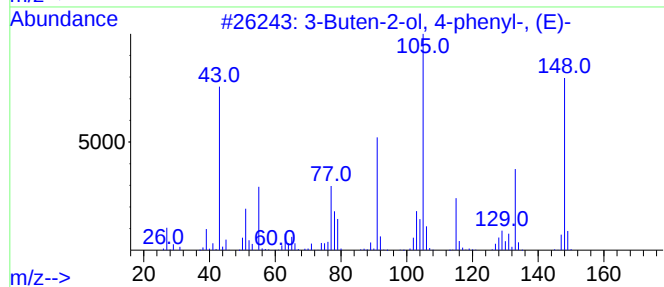
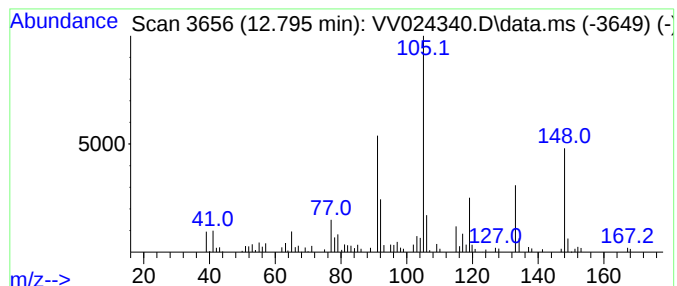
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Quant Title : VOC Analysis

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

Peak Number 28 3-Buten-2-ol, 4-phenyl-, (E)- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.795	3.09 ug/L	57566	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	50
2			2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	47
3			2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	43
4			Benzene, (1-ethylpropyl)-	148	C11H16	001196-58-3	43
5			Benzene, (1,2-dimethylpropyl)-	148	C11H16	004481-30-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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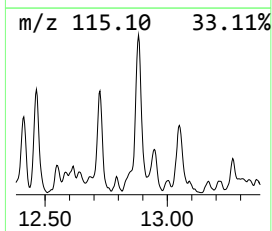
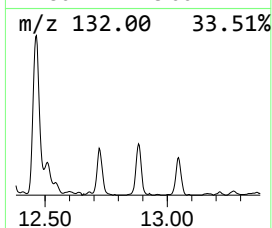
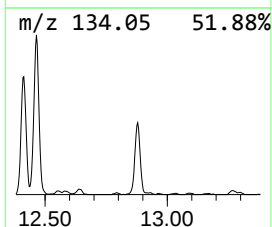
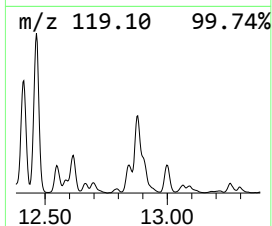
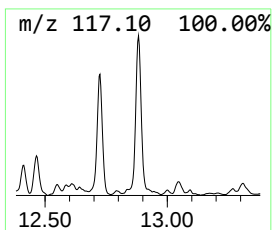
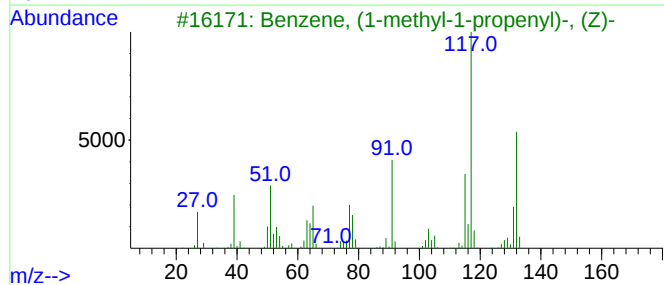
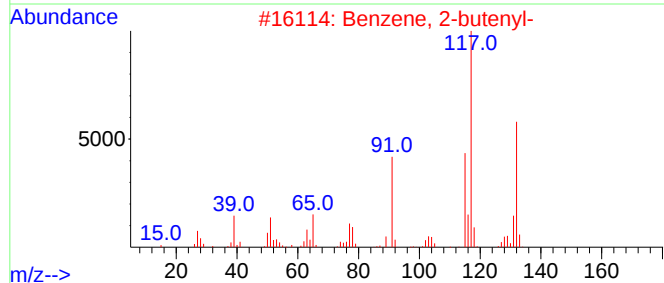
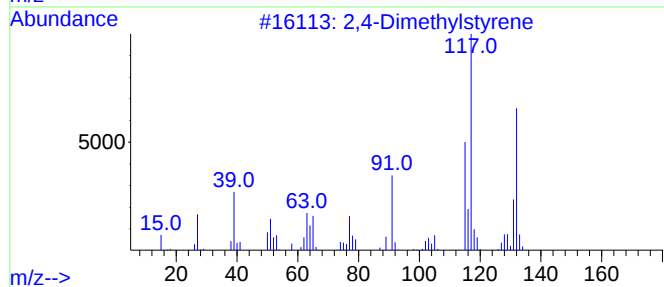
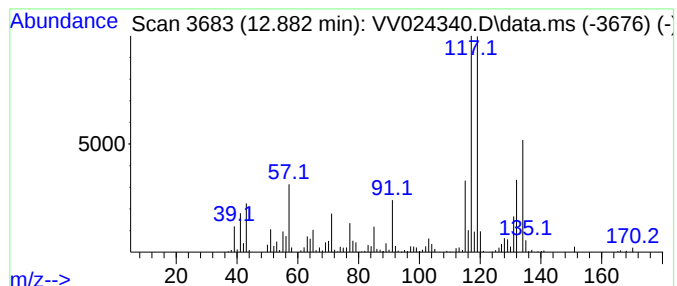
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Quant Title : VOC Analysis

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

Peak Number 29 2,4-Dimethylstyrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	20.03 ug/L	373122	1,4-Dichlorobenzene-d4	11.249

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,4-Dimethylstyrene	132	C10H12	002234-20-0	78
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	50
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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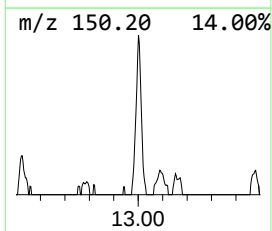
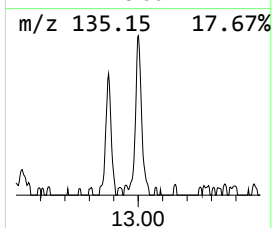
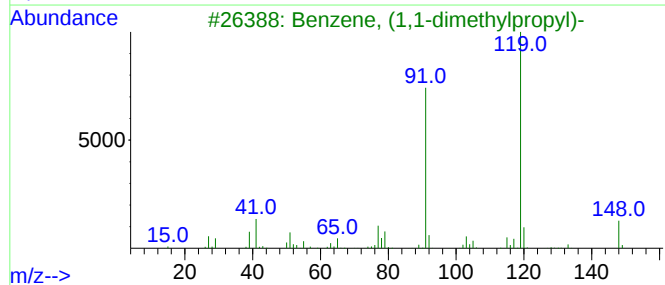
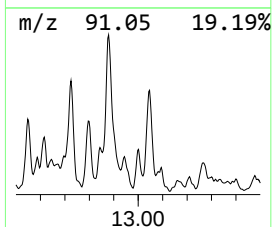
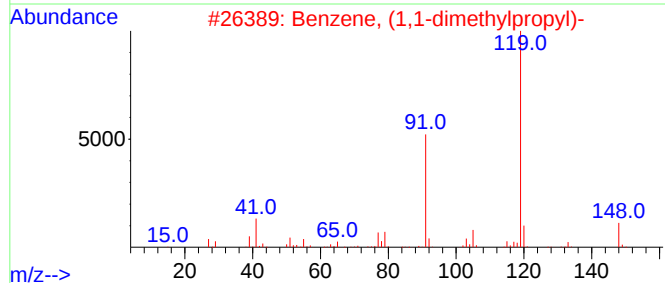
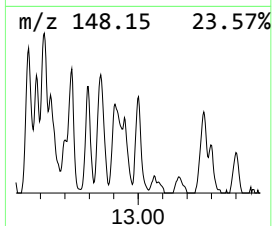
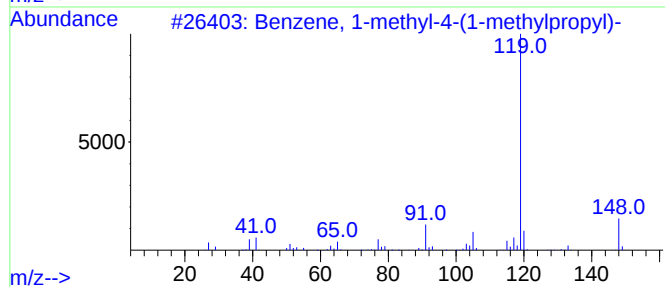
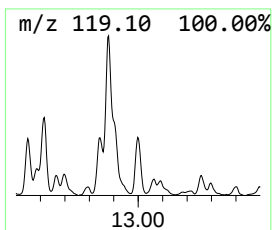
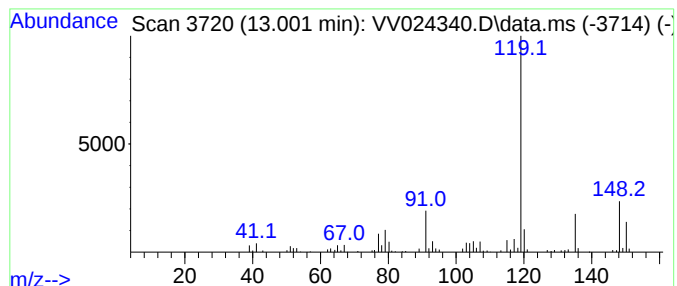
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Quant Title : VOC Analysis

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

Peak Number 30 Benzene, 1-methyl-4-(1-meth... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.001	2.53 ug/L	47156	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	72
2			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	59
3			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	59
4			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	59
5			p-Cymene	134	C10H14	000099-87-6	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

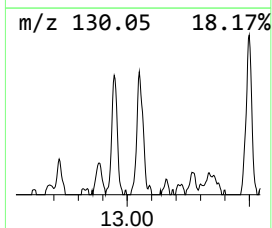
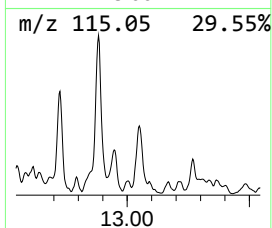
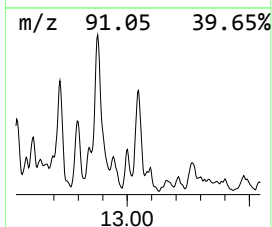
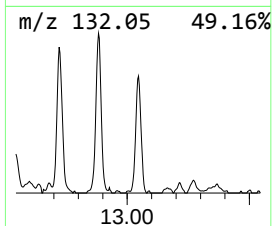
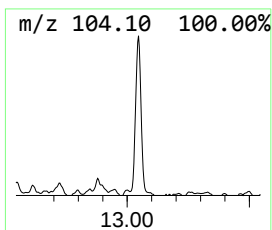
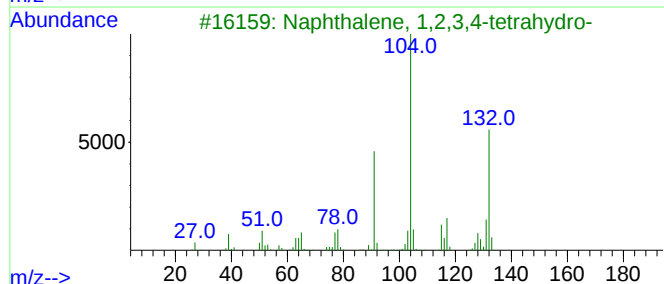
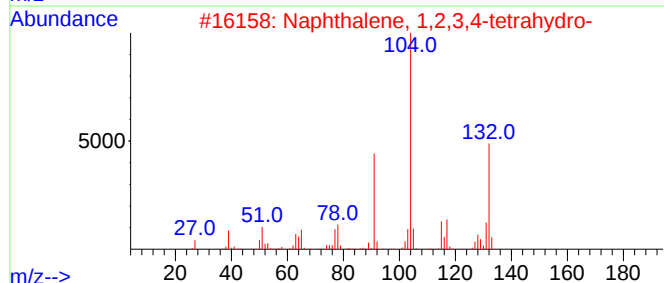
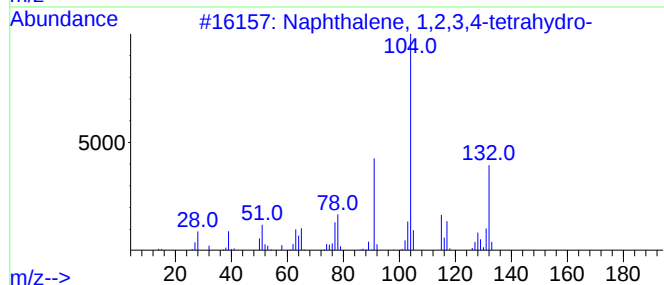
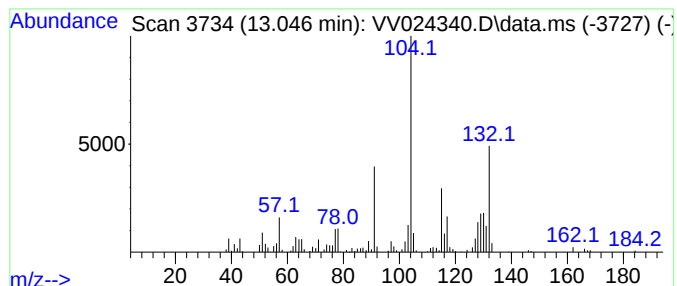
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Quant Title : VOC Analysis

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

Peak Number 31 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.046	6.78 ug/L	126367	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
3			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
4			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	81
5			Indan, epoxide	132	C9H8O	000768-22-9	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

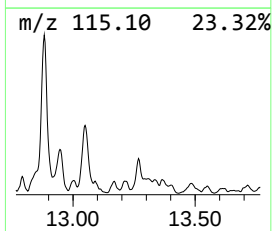
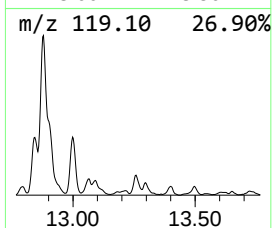
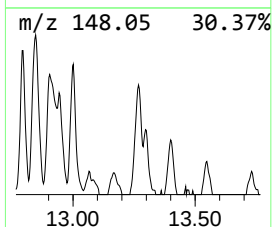
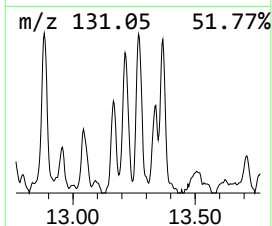
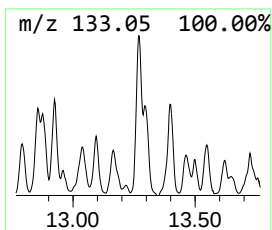
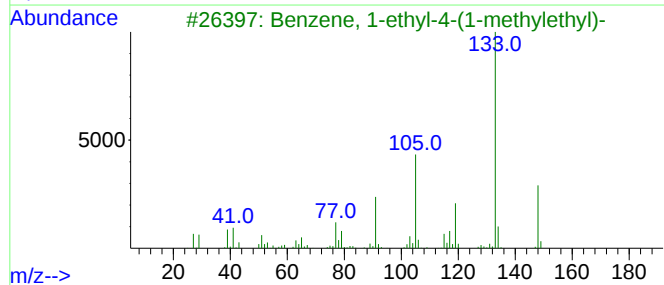
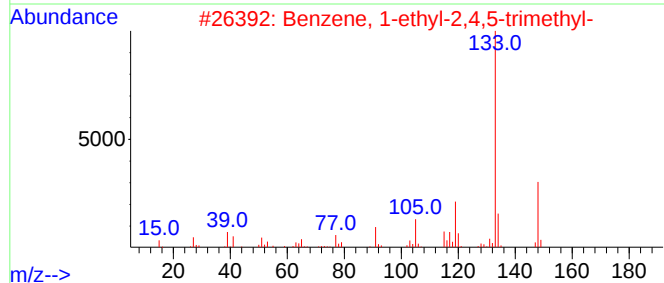
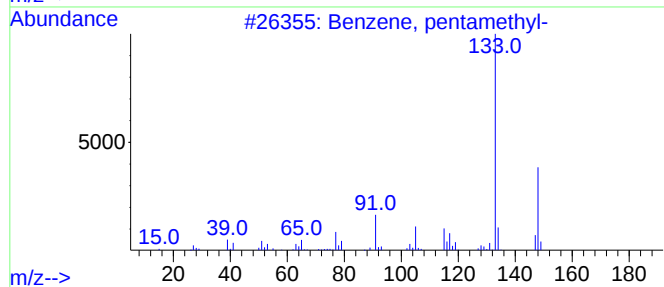
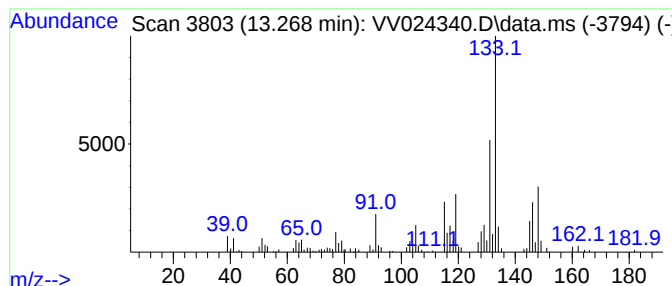
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Quant Title : VOC Analysis

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

Peak Number 32 Benzene, pentamethyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	4.53 ug/L	84381	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	55
2			Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	49
3			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	49
4			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	49
5			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

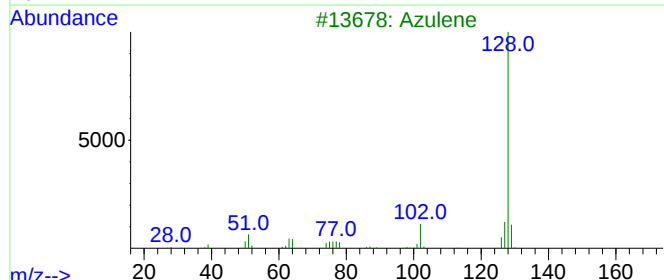
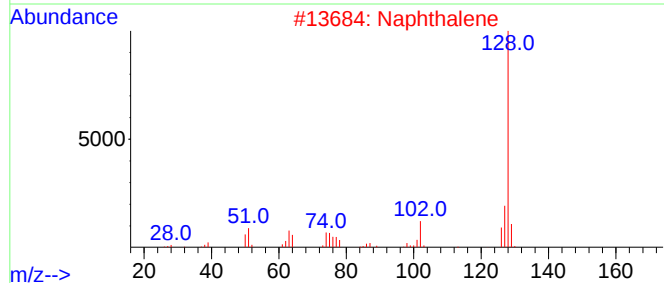
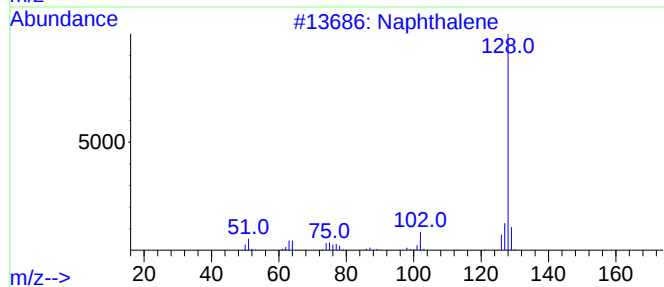
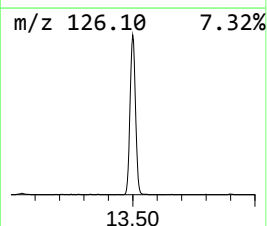
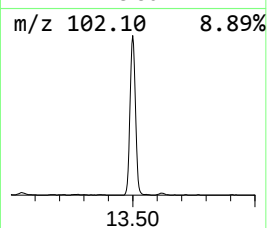
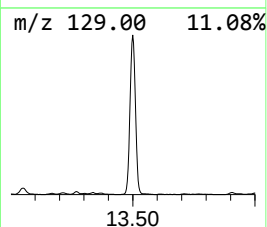
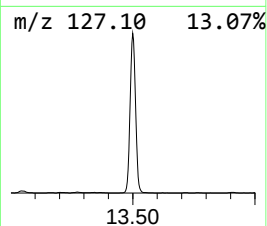
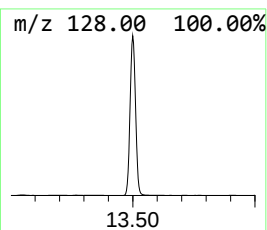
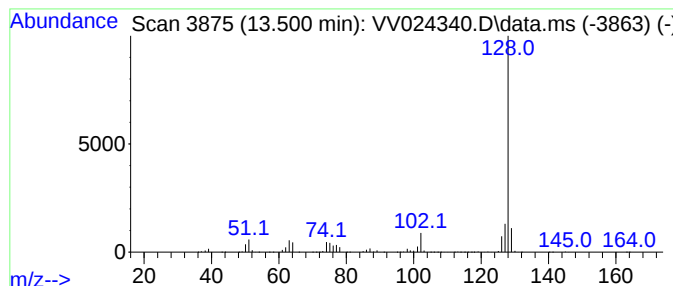
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Quant Title : VOC Analysis

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

Peak Number 33 Azulene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	112.03 ug/L	2086910	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Azulene	128	C10H8	000275-51-4	93
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

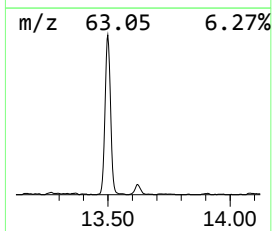
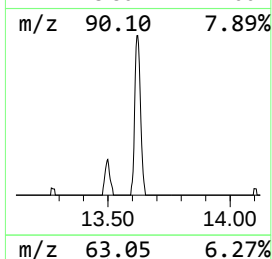
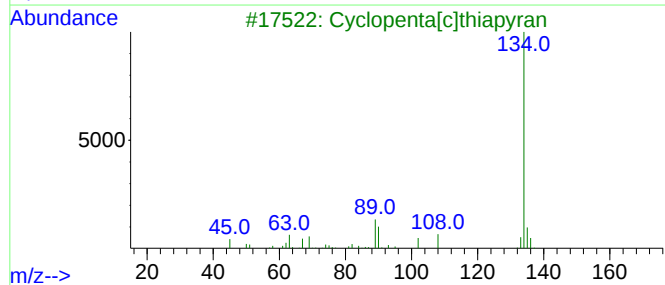
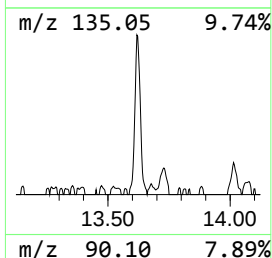
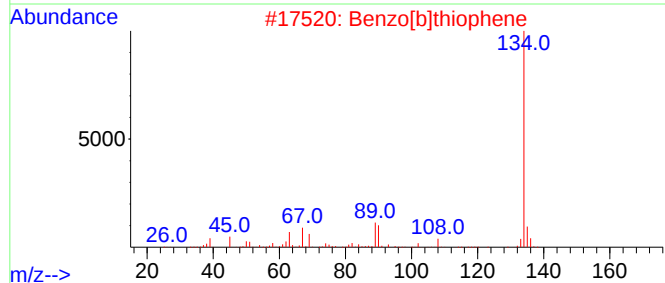
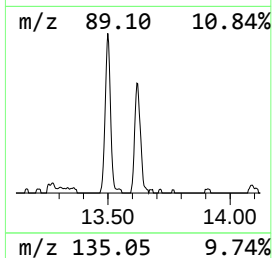
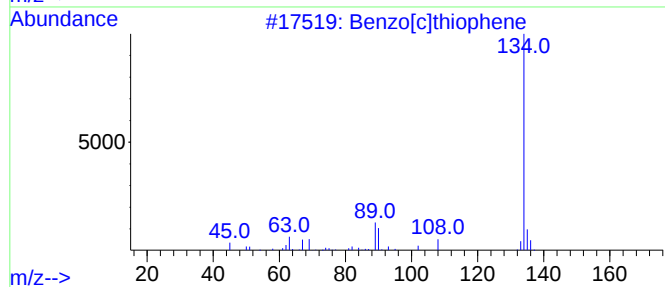
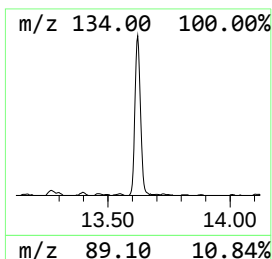
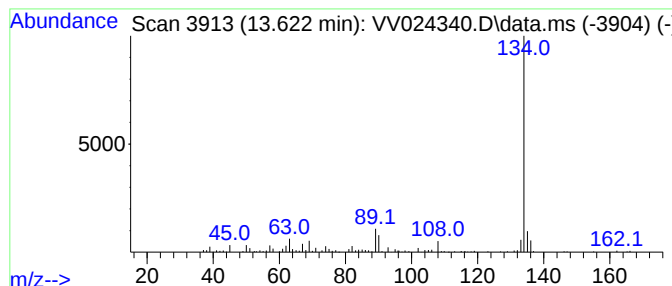
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Quant Title : VOC Analysis

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

Peak Number 34 Benzo[c]thiophene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.622	7.78 ug/L	144856	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]thiophene	134	C8H6S	000270-82-6	97
2			Benzo[b]thiophene	134	C8H6S	000095-15-8	94
3			Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
4			Benzo[b]thiophene	134	C8H6S	000095-15-8	90
5			Benzo[b]thiophene	134	C8H6S	000095-15-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

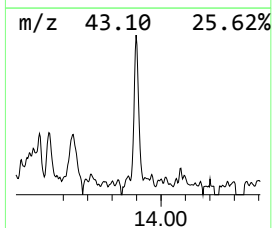
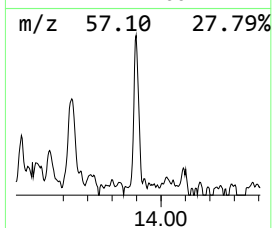
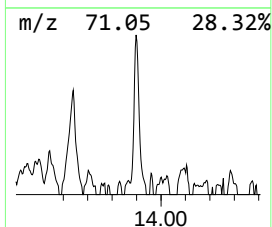
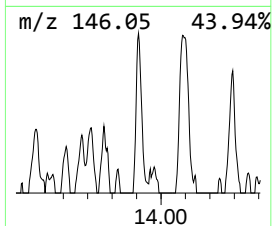
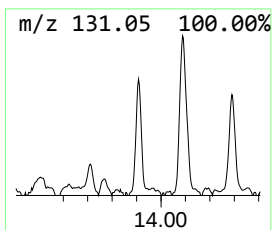
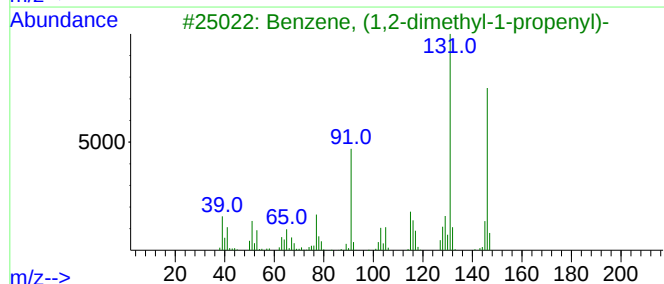
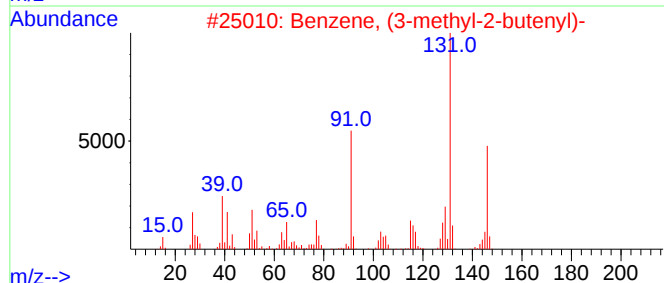
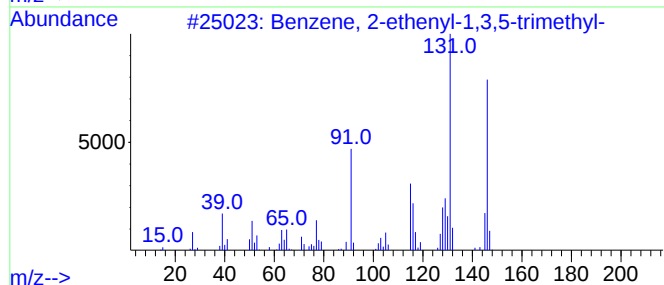
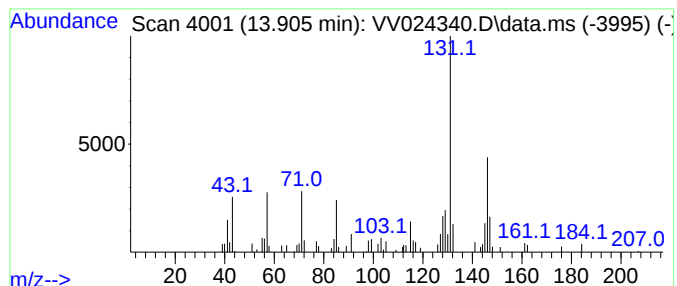
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Quant Title : VOC Analysis

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

Peak Number 35 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.905	2.59 ug/L	48160	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	76
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	76
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	76
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	70
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

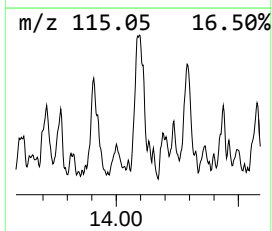
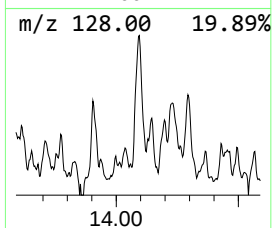
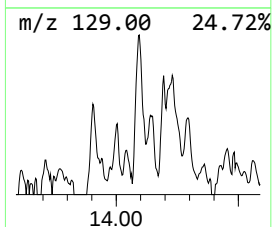
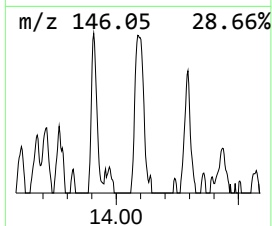
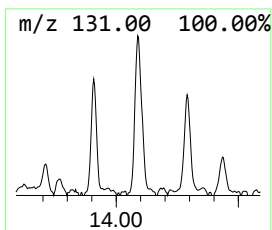
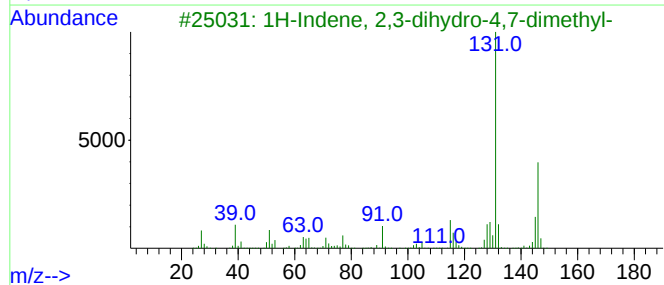
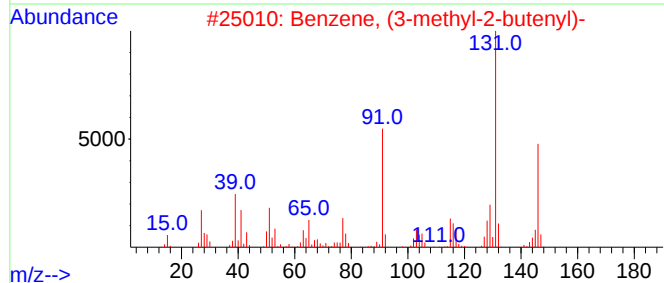
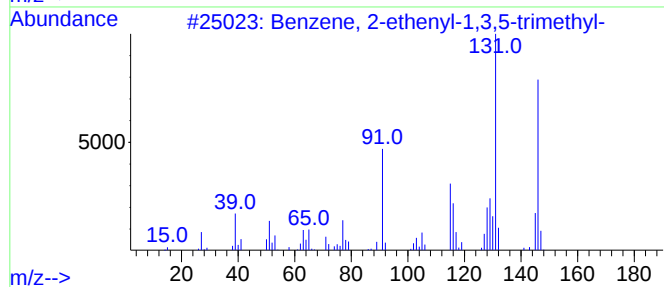
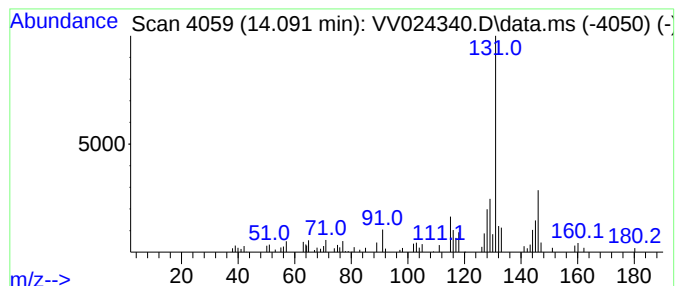
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Quant Title : VOC Analysis

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

Peak Number 36 Benzene, (3-methyl-2-butenyl)- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.091	2.95 ug/L	54865	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	81
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	76
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

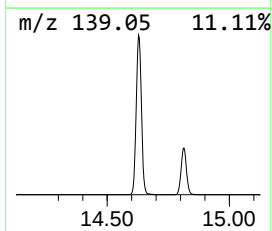
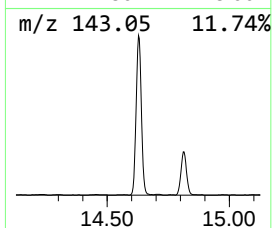
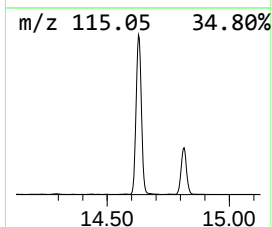
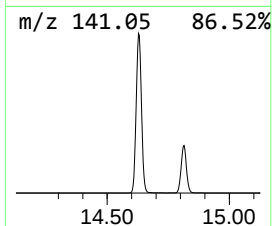
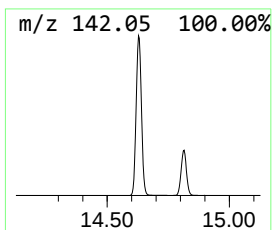
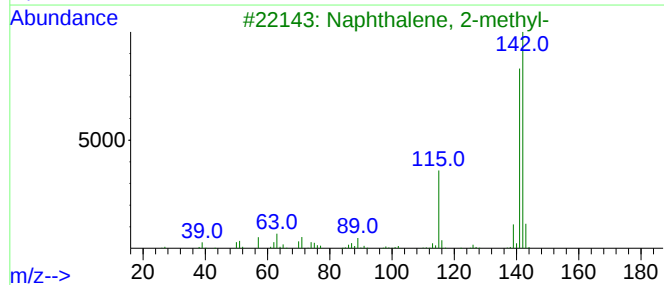
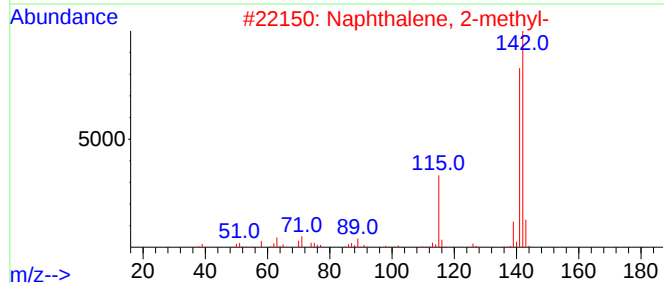
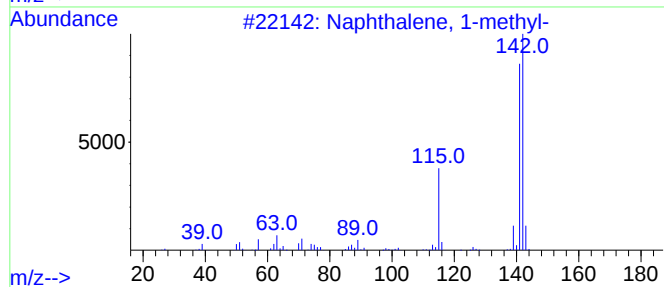
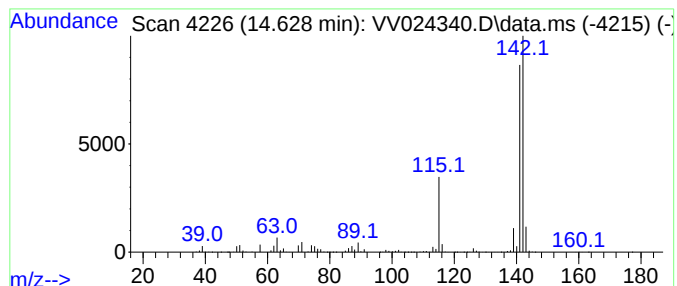
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Quant Title : VOC Analysis

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

Peak Number 37 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.628	152.26 ug/L	2836320	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

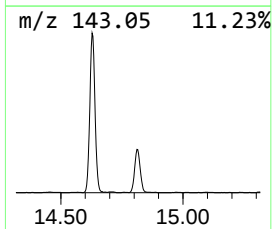
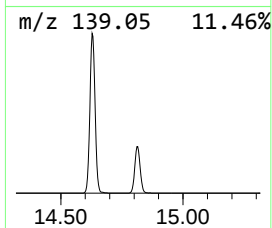
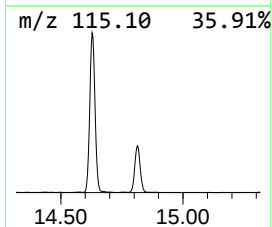
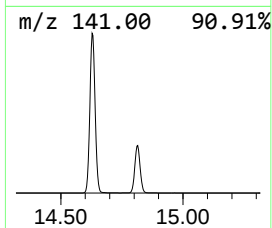
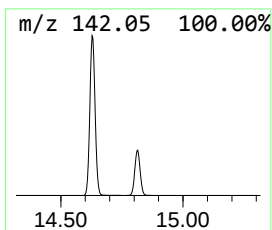
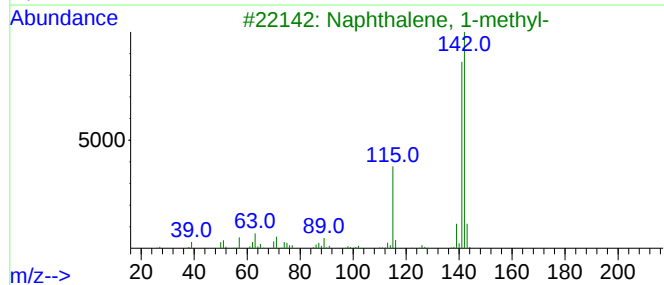
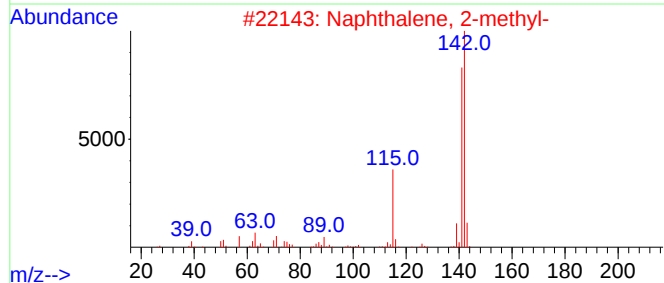
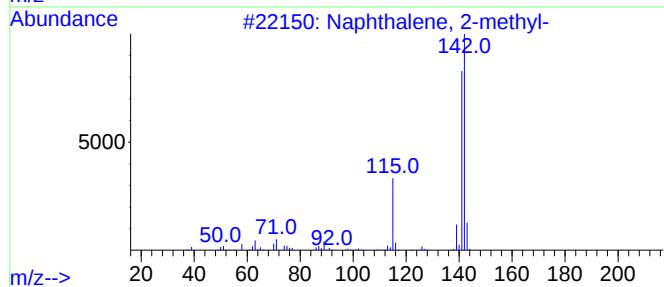
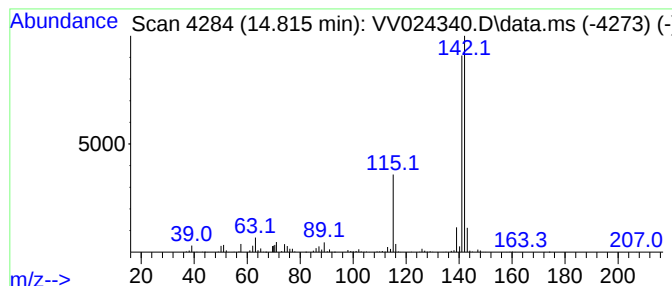
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Quant Title : VOC Analysis

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

Peak Number 38 Naphthalene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.815	45.44 ug/L	846383	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

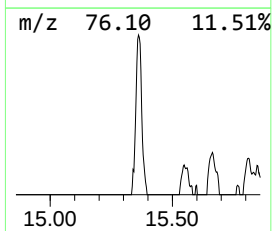
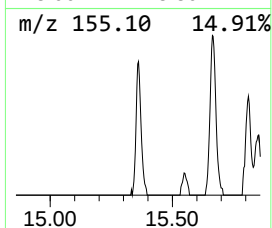
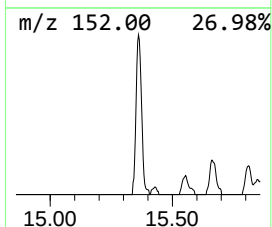
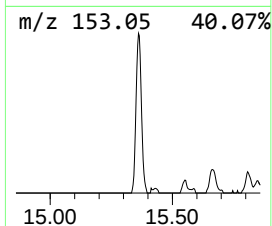
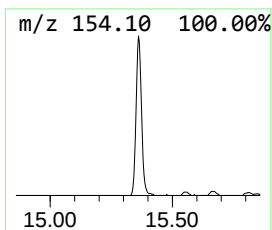
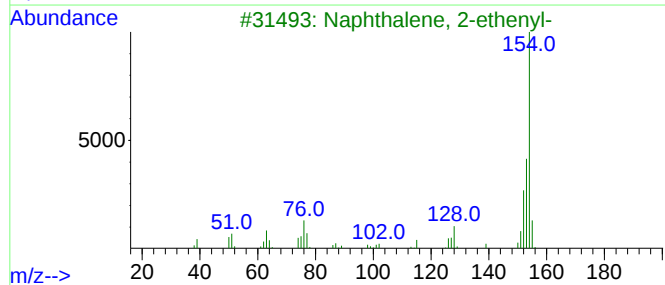
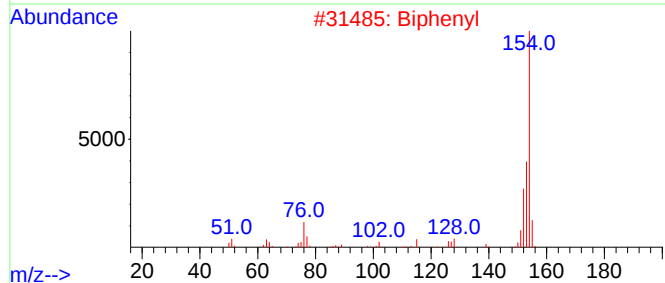
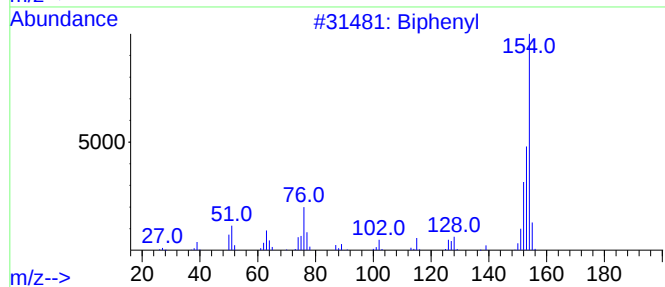
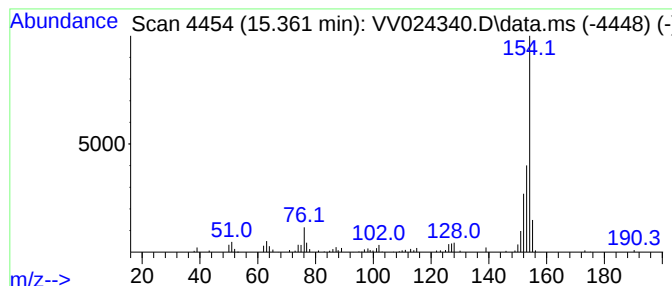
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Quant Title : VOC Analysis

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

Peak Number 39 Biphenyl Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.361	4.21 ug/L	78422	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenyl	154	C12H10	000092-52-4	93
2			Biphenyl	154	C12H10	000092-52-4	93
3			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	91
4			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	91
5			Biphenyl	154	C12H10	000092-52-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV011222\
Data File : VV024340.D
Acq On : 12 Jan 2022 19:10
Operator : SY/MD
Sample : N1084-09 20X
Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BGLN2

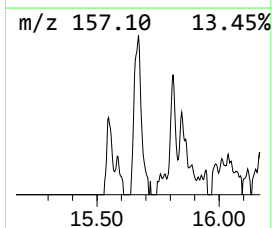
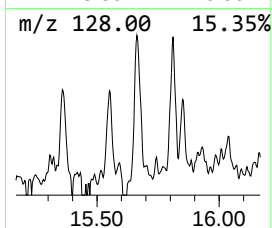
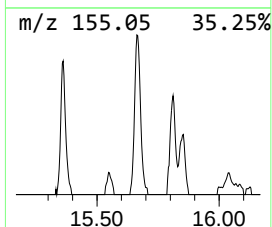
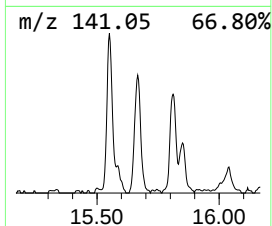
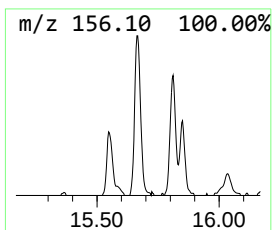
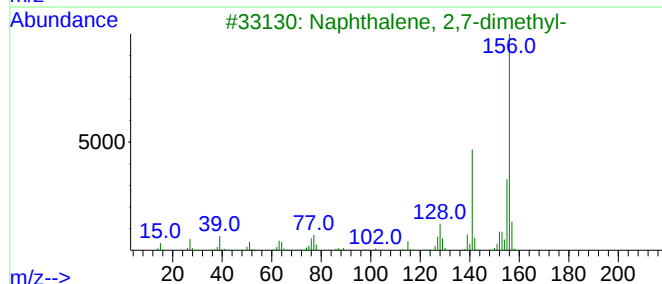
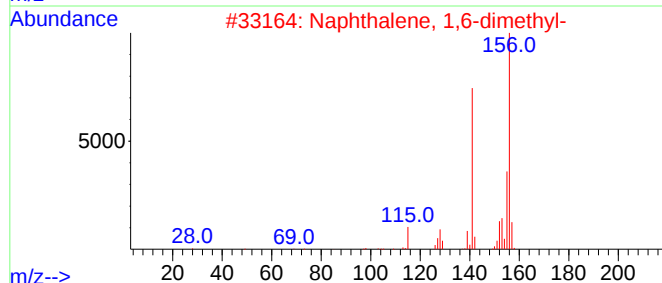
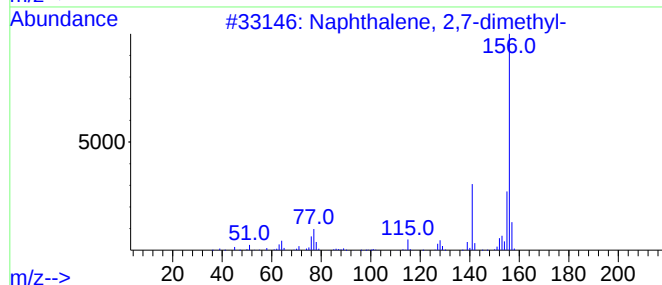
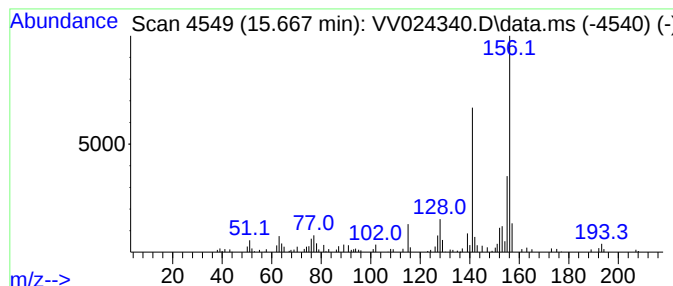
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Quant Title : VOC Analysis

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

Peak Number 40 Naphthalene, 2,7-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.667	4.07 ug/L	75881	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW011222\
 Data File : VW024340.D
 Acq On : 12 Jan 2022 19:10
 Operator : SY/MD
 Sample : N1084-09 20X
 Misc : 6.39g/5.0mL/100ul/5.0mL/MSVOA_V/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BGLN2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM123121WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	8.352	2.7	ug/L	39166	2	8.850	718182	50.0
(DEL) Alkane: S...	8.667	6.4	ug/L	92449	2	8.850	718182	50.0
(DEL) Alkane: S...	8.789	6.1	ug/L	87925	2	8.850	718182	50.0
(DEL) Alkane: S...	9.204	35.0	ug/L	502441	2	8.850	718182	50.0
(DEL) Alkane: C...	9.316	2.7	ug/L	39203	2	8.850	718182	50.0
unknown-01	9.474	9.7	ug/L	139252	2	8.850	718182	50.0
(DEL) Alkane: C...	9.673	3.9	ug/L	56093	2	8.850	718182	50.0
(DEL) Alkane: S...	9.709	24.1	ug/L	346295	2	8.850	718182	50.0
(DEL) Alkane: C...	9.799	20.1	ug/L	288880	2	8.850	718182	50.0
(DEL) Alkane: S...	10.017	9.8	ug/L	140992	2	8.850	718182	50.0
Benzene, propyl-	10.352	32.6	ug/L	607309	3	11.249	931419	50.0
Benzene, 1-ethy...	10.445	161.1	ug/L	3000180	3	11.249	931419	50.0
Benzene, 1-ethy...	10.734	57.2	ug/L	1065950	3	11.249	931419	50.0
(DEL) Alkane: S...	11.034	8.6	ug/L	159293	3	11.249	931419	50.0
(DEL) Alkane: C...	11.152	34.4	ug/L	640398	3	11.249	931419	50.0
Benzene, 1,2,3-...	11.336	68.8	ug/L	1280660	3	11.249	931419	50.0
(DEL) Alkane: S...	11.461	13.4	ug/L	248818	3	11.249	931419	50.0
Indane	11.525	253.8	ug/L	4728670	3	11.249	931419	50.0
Benzene, 2-ethy...	11.911	17.6	ug/L	327390	3	11.249	931419	50.0
Benzene, 2-ethy...	12.014	34.3	ug/L	639559	3	11.249	931419	50.0
1-Methyl-2-phen...	12.117	7.6	ug/L	140970	3	11.249	931419	50.0
Benzene, 1,3-di...	12.255	13.2	ug/L	246701	3	11.249	931419	50.0
(DEL) Alkane: C...	12.368	15.3	ug/L	284728	3	11.249	931419	50.0
Benzene, 1,2,4,...	12.413	13.5	ug/L	252087	3	11.249	931419	50.0
Benzene, 1,3-di...	12.464	37.0	ug/L	688778	3	11.249	931419	50.0
Benzene, 1-meth...	12.725	10.5	ug/L	194836	3	11.249	931419	50.0
3-Buten-2-ol, 4...	12.795	3.1	ug/L	57566	3	11.249	931419	50.0
2,4-Dimethylsty...	12.882	20.0	ug/L	373122	3	11.249	931419	50.0
Benzene, 1-meth...	13.001	2.5	ug/L	47156	3	11.249	931419	50.0
Naphthalene, 1,...	13.046	6.8	ug/L	126367	3	11.249	931419	50.0
Benzene, pentam...	13.268	4.5	ug/L	84381	3	11.249	931419	50.0
Azulene	13.500	112.0	ug/L	2086910	3	11.249	931419	50.0
Benzo[c]thiophene	13.622	7.8	ug/L	144856	3	11.249	931419	50.0
Benzene, 2-ethe...	13.905	2.6	ug/L	48160	3	11.249	931419	50.0
Benzene, (3-met...	14.091	3.0	ug/L	54865	3	11.249	931419	50.0
Naphthalene, 1-...	14.628	152.3	ug/L	2836320	3	11.249	931419	50.0
Naphthalene, 2-...	14.815	45.4	ug/L	846383	3	11.249	931419	50.0
Biphenyl	15.361	4.2	ug/L	78422	3	11.249	931419	50.0
Naphthalene, 2,...	15.667	4.1	ug/L	75881	3	11.249	931419	50.0