Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Title : SW846 8260

Signal : TIC: VX032639.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.301	28	35	40	rBV5	4743	14058	2.93%	0.203%
2	2.380	205	212	221	rBV	50489	80001	16.69%	1.154%
3	2.532	229	237	257	rBV	265995	478780	99.86%	6.908%
4	4.574	564	572	577	rBV3	1802	4814	1.00%	0.069%
5	5.032	637	647	657	rBB3	1601	4952	1.03%	0.071%
6	5.398	697	707	717	rBV2	22104	65049	13.57%	0.938%
7	5.562	722	734	749	rBV	93298	259267	54.08%	3.741%
8	5.964	790	800	810	rBV2	23401	62194	12.97%	0.897%
9	6.769	922	932	949	rBV	119617	290347	60.56%	4.189%
10	8.653	1234	1241	1248	rBV	250664	380932	79.45%	5.496%
11	8.726	1248	1253	1260	rVB	12252	19957	4.16%	0.288%
12	10.055	1466	1471	1482	rBV	194708	258027	53.82%	3.723%
13	10.195	1490	1494	1502	rBV	20780	28381	5.92%	0.409%
14	10.299	1506	1511	1519	rBV	117382	164993	34.41%	2.380%
15	10.647	1561	1568	1574	rBV	78131	103925	21.68%	1.499%
16	10.964	1615	1620	1626	rVB	8366	9987	2.08%	0.144%
17	11.086	1634	1640	1645	rBV	148007	195720	40.82%	2.824%
18	11.305	1671	1676	1683	rBV	22100	27257	5.69%	0.393%
19	11.378	1683	1688	1696	rVV	168484	277850	57.95%	4.009%
20	11.451	1696	1700	1711	rVB	95823	116394	24.28%	1.679%
21	11.616	1718	1727	1736	rBV	84873	109319	22.80%	1.577%
22	11.750	1744	1749	1757	rVB	389604	479444	100.00%	6.917%
23	11.890	1769	1772	1776	rVB	7217	8453	1.76%	0.122%
24	11.969	1781	1785	1789	rVV	32019	37732	7.87%	0.544%
25	12.024	1789	1794	1799	rVV	177460	228366	47.63%	3.295%
26	12.085	1799	1804	1811	rVB	158836	197921	41.28%	2.855%
27	12.244	1823	1830	1833	rBV	129478	184579	38.50%	2.663%
28	12.317	1838	1842	1852	rVB2	89244	127244	26.54%	1.836%
29	12.408	1852	1857	1861	rBV2	10258	14621	3.05%	0.211%
30	12.463	1861	1866	1871	rVB	31512	35767	7.46%	0.516%
31	12.530	1872	1877	1879	rBV	66463	79982	16.68%	1.154%
32	12.555	1879	1881	1886	rVB	62644	65503	13.66%	0.945%
33	12.616	1886	1891	1894	rBV	92548	115972	24.19%	1.673%
34	12.646	1894	1896	1900	rVB	28823	29865	6.23%	0.431%
35	12.701	1900	1905	1914	rBV2	82894	122964	25.65%	1.774%
36	12.799	1917	1921	1924	rBV4	4469	5235	1.09%	0.076%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Title : SW846 8260

37	12.847	1924	1929	1935	rBV2	39017	50627	10.56%	0.730%
38	12.921	1935	1941	1944	rBV	64035	76786	16.02%	1.108%
39	12.963	1944	1948	1954	rVB	106836	125515	26.18%	1.811%
40	13.024	1954	1958	1960	rBV2	10083	13514	2.82%	0.195%
41	13.140	1973	1977	1978	rVV2	10990	13363	2.79%	0.193%
42	13.170	1978	1982	1986	rVV	99511	128094	26.72%	1.848%
43	13.207	1986	1988	1992	rVV3	8320	10610	2.21%	0.153%
44	13.256	1992	1996	1997	rVV2	12911	16612	3.46%	0.240%
45	13.292	1997	2002	2007	rVV2	214561	295781	61.69%	4.267%
46	13.341	2007	2010	2013	rVV4	11132	18359	3.83%	0.265%
47	13.372	2013	2015	2018	rVV	9121	9769	2.04%	0.141%
48	13.420	2018	2023	2031	rVB	105921	133849	27.92%	1.931%
49	13.506	2031	2037	2040	rBV2	17373	24078	5.02%	0.347%
50	13.542	2040	2043	2046	rVV	37030	49365	10.30%	0.712%
51	13.585	2046	2050	2055	rVV3	39524	74023	15.44%	1.068%
52	13.664	2055	2063	2070	rVB2	40700	85619	17.86%	1.235%
53	13.774	2075	2081	2089	rBV	115334	157034	32.75%	2.266%
54	13.853	2089	2094	2100	rVB	21003	27325	5.70%	0.394%
55	13.926	2100	2106	2110	rBV2	32819	45963	9.59%	0.663%
56	13.969	2110	2113	2117	rVB	16821	19555	4.08%	0.282%
57	14.073	2126	2130	2137	rBV	38007	49250	10.27%	0.711%
58	14.225	2149	2155	2163	rBV3	55005	111834	23.33%	1.613%
59	14.323	2167	2171	2176	rBV	18173	25699	5.36%	0.371%
60	14.371	2176	2179	2184	rVB	33544	38941	8.12%	0.562%
61	14.481	2191	2197	2202	rBV2	30526	40475	8.44%	0.584%
62	14.634	2210	2222	2227	rBV	273572	355536	74.16%	5.129%
63	14.780	2240	2246	2251	rBV	143892	181632	37.88%	2.620%
64	15.182	2309	2312	2319	rVB3	5225	7975	1.66%	0.115%
65	15.377	2339	2344	2348	rBV	6769	9858	2.06%	0.142%
66	15.475	2354	2360	2371	rBV	11986	19887	4.15%	0.287%
67	15.615	2377	2383	2386	rBV	13465	20015	4.17%	0.289%
68	15.652	2386	2389	2394	rVB2	5976	8470	1.77%	0.122%

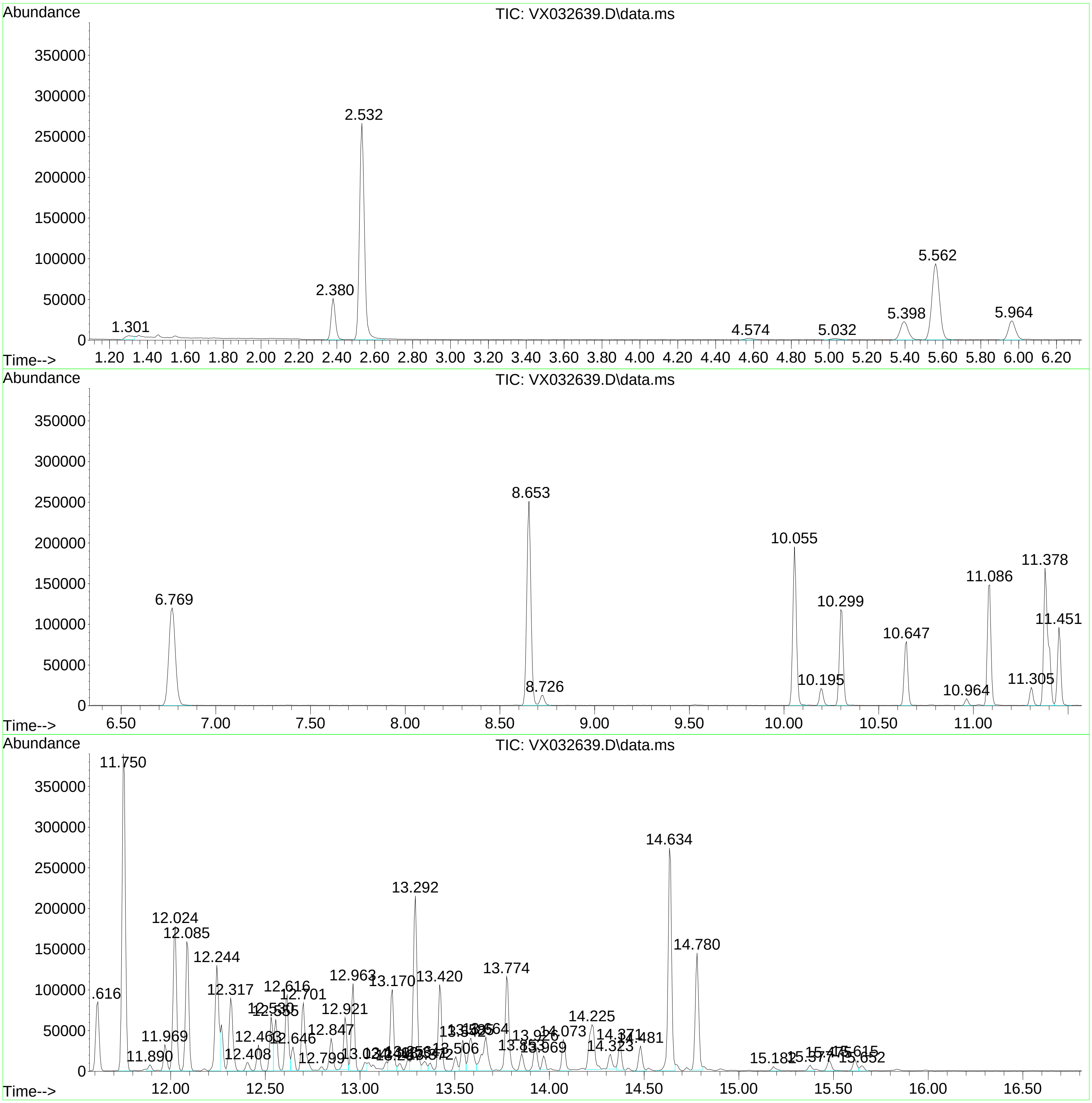
Sum of corrected areas: 6931265

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
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Quant Title : SW846 8260

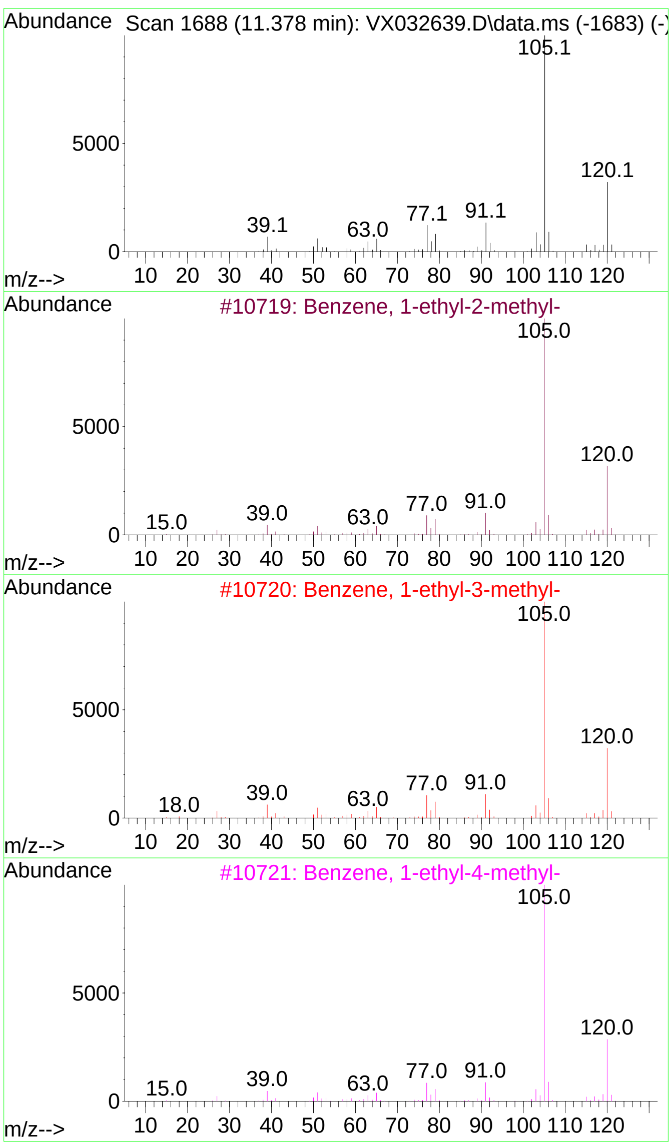
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

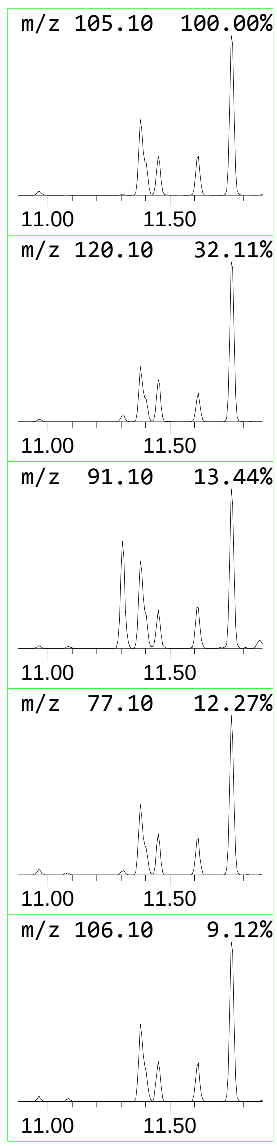
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	60.83 ug/l	277850	1,4-Dichlorobenzene-d4	12.024

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-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

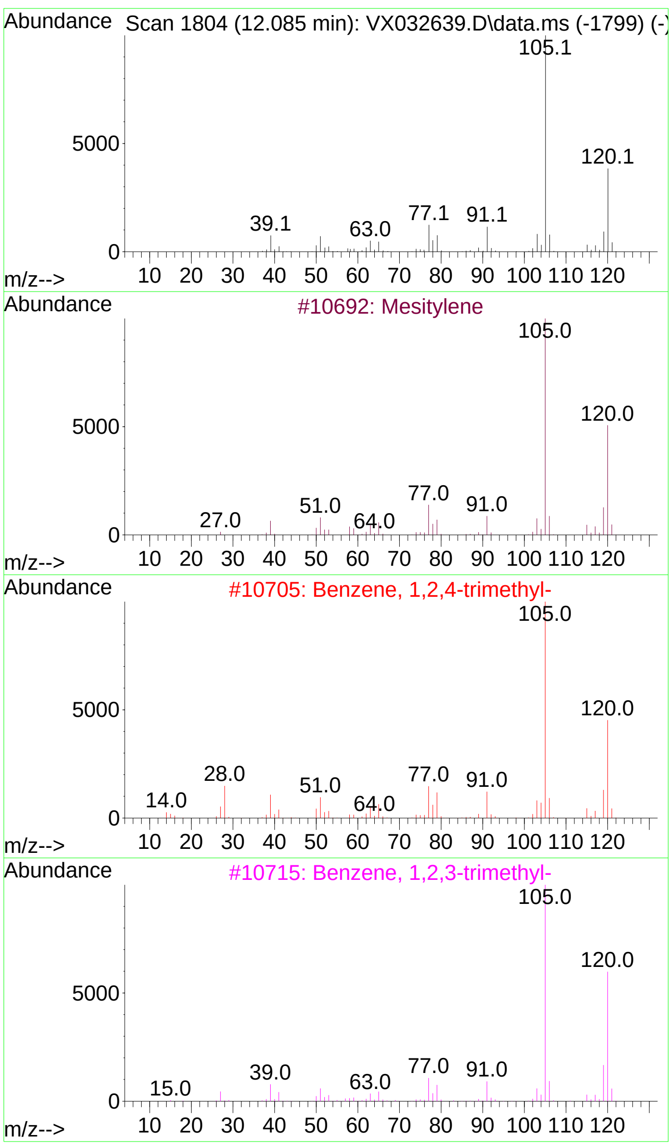
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

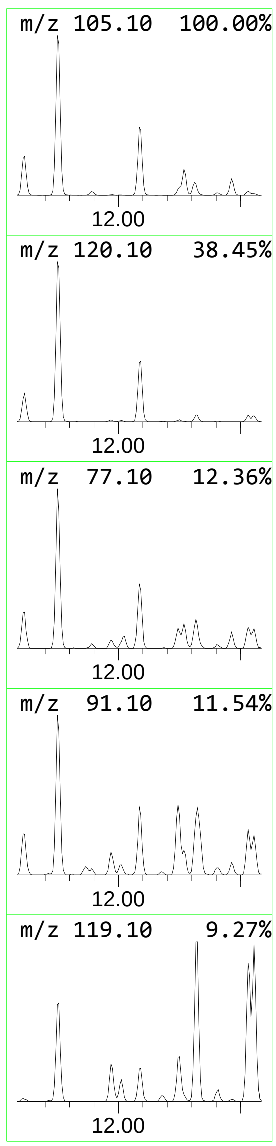
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	43.33 ug/l	197921	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

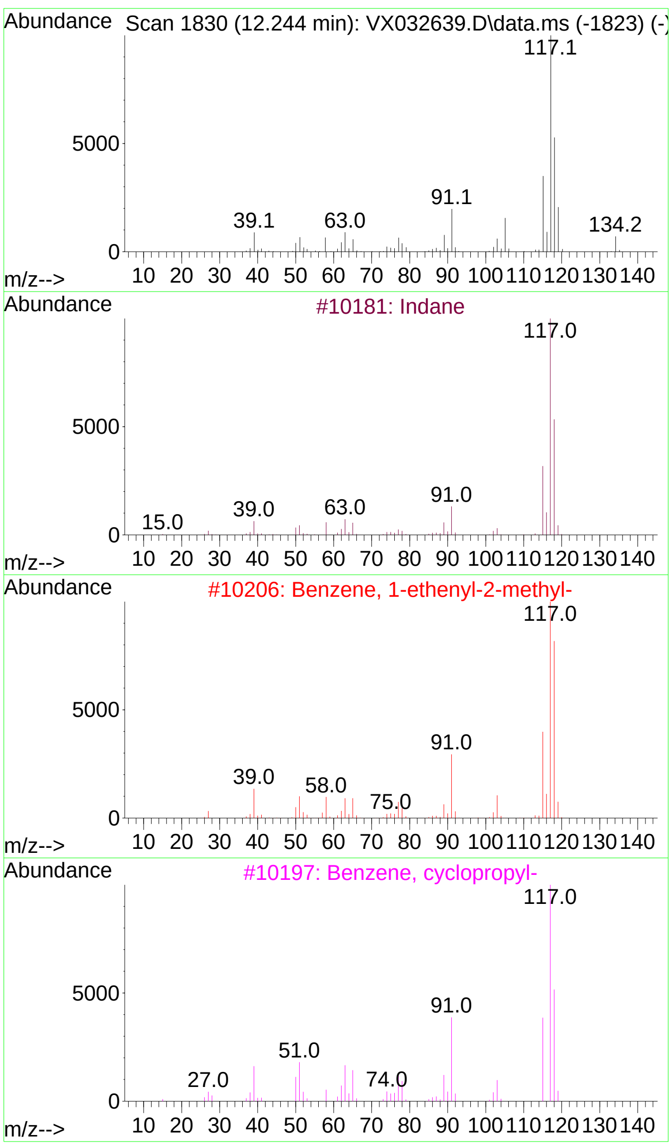
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

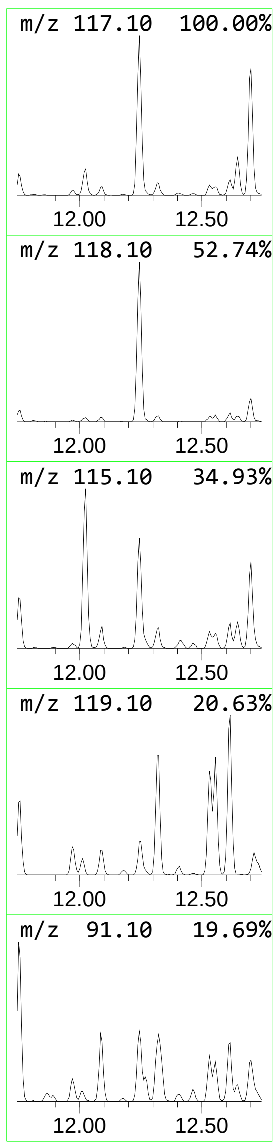
****

Peak Number 3 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	40.41 ug/l	184579	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
5	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

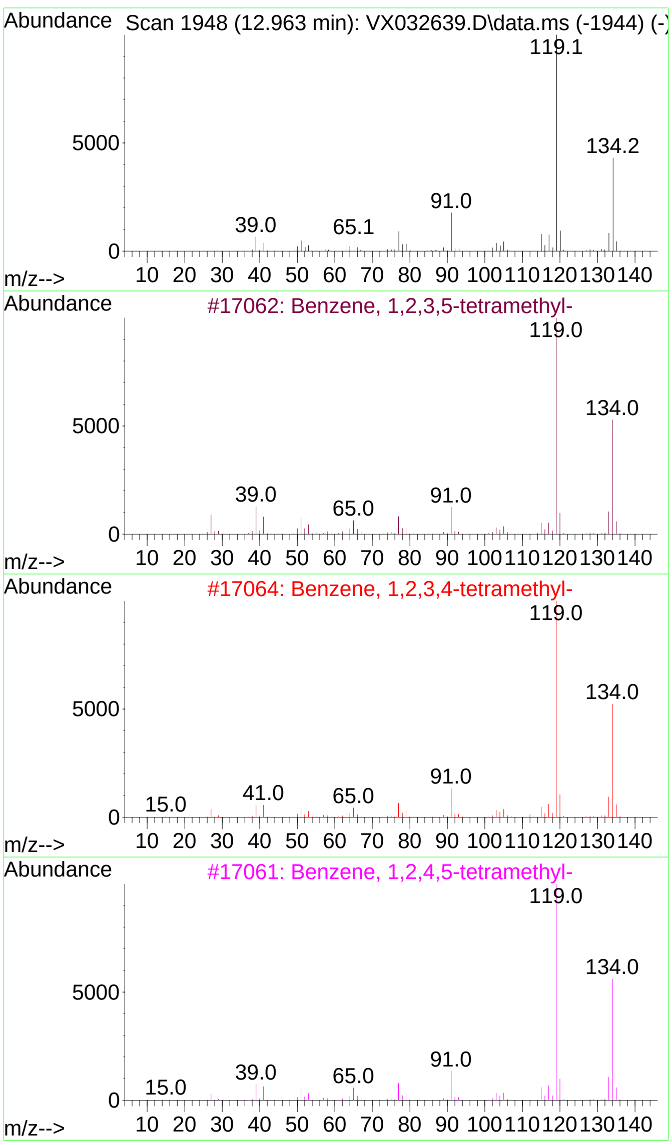
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

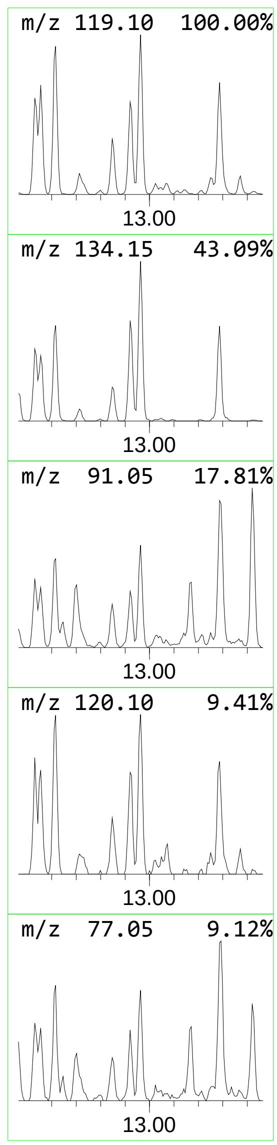
****

Peak Number 4 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	27.48 ug/l	125515	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





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Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On    : 07 Nov 2022  19:51
Operator  : JC/MD
Sample    : N5504-01
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 26   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title  : SW846 8260

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

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

```
**********************************************************************
Peak Number  5  Benzene, 1-ethenyl-4-ethyl-     Concentration Rank  9

  R.T.     EstConc          Area       Relative to ISTD         R.T.
----------------------------------------------------------------------
 13.171   28.05 ug/l       128094   1,4-Dichlorobenzene-d4     12.024

 Hit# of  5    Tentative ID              MW  MolForm       CAS#      Qual
----------------------------------------------------------------------
 1 Benzene, 1-ethenyl-4-ethyl-         132 C10H12        003454-07-7 94
 2 Benzene, 2-ethenyl-1,4-dimethyl-    132 C10H12        002039-89-6 93
 3 1H-Indene, 2,3-dihydro-5-methyl-    132 C10H12        000874-35-1 93
 4 Benzene, (1-methyl-1-propenyl)-,... 132 C10H12        000768-00-3 91
 5 1H-Indene, 2,3-dihydro-4-methyl-    132 C10H12        000824-22-6 90
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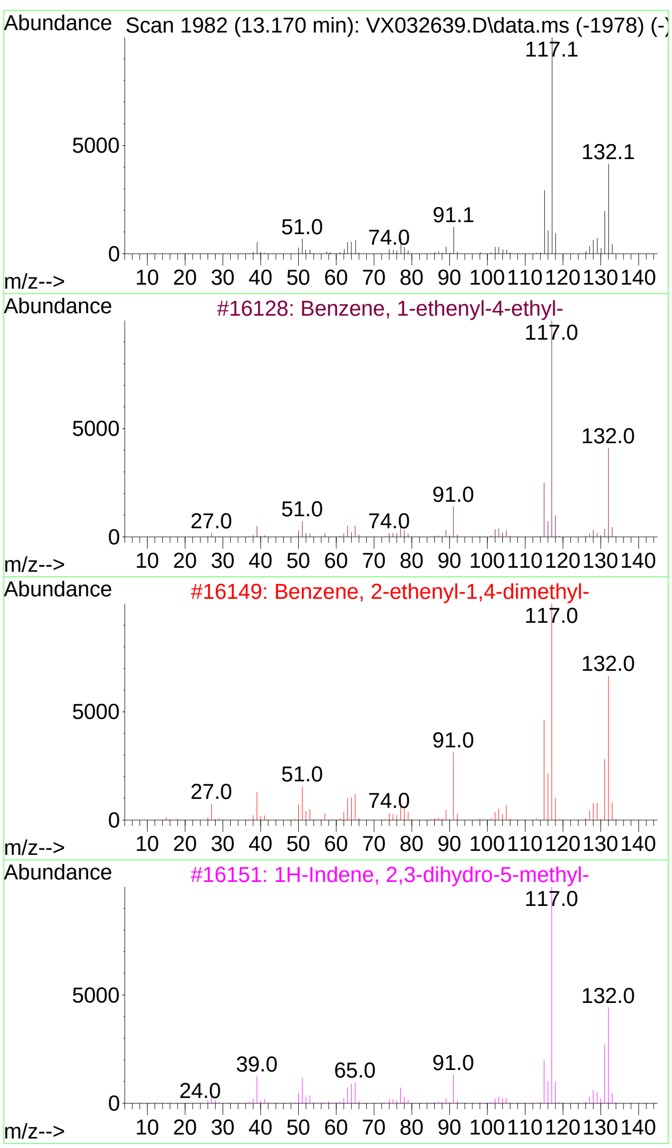

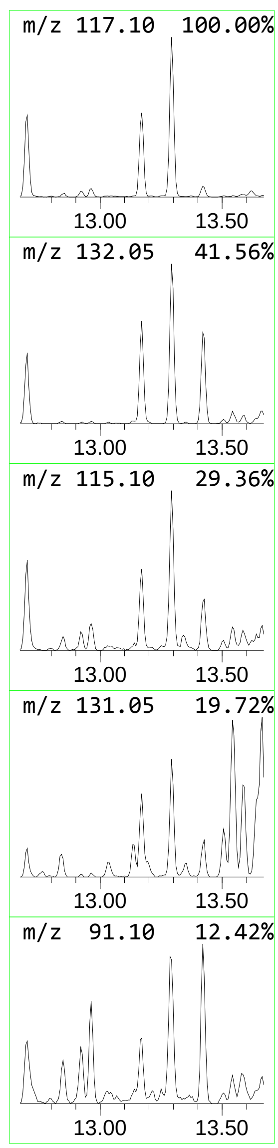

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Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On    : 07 Nov 2022  19:51
Operator  : JC/MD
Sample    : N5504-01
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 26   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title  : SW846 8260

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

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

```
*********************************************************************
Peak Number  6  1-Phenyl-1-butene              Concentration Rank  2

  R.T.     EstConc          Area       Relative to ISTD         R.T.
---------------------------------------------------------------------
 13.292   64.76 ug/l       295781   1,4-Dichlorobenzene-d4     12.024

 Hit# of  5     Tentative ID              MW  MolForm       CAS#      Qual
---------------------------------------------------------------------
 1 1-Phenyl-1-butene                     132 C10H12        000824-90-8 81
 2 3-Phenylbut-1-ene                     132 C10H12        000934-10-1 81
 3 Benzene, 1-ethenyl-3-ethyl-           132 C10H12        007525-62-4 76
 4 Benzene, 1-ethenyl-4-ethyl-           132 C10H12        003454-07-7 76
 5 Indan, 1-methyl-                      132 C10H12        000767-58-8 76
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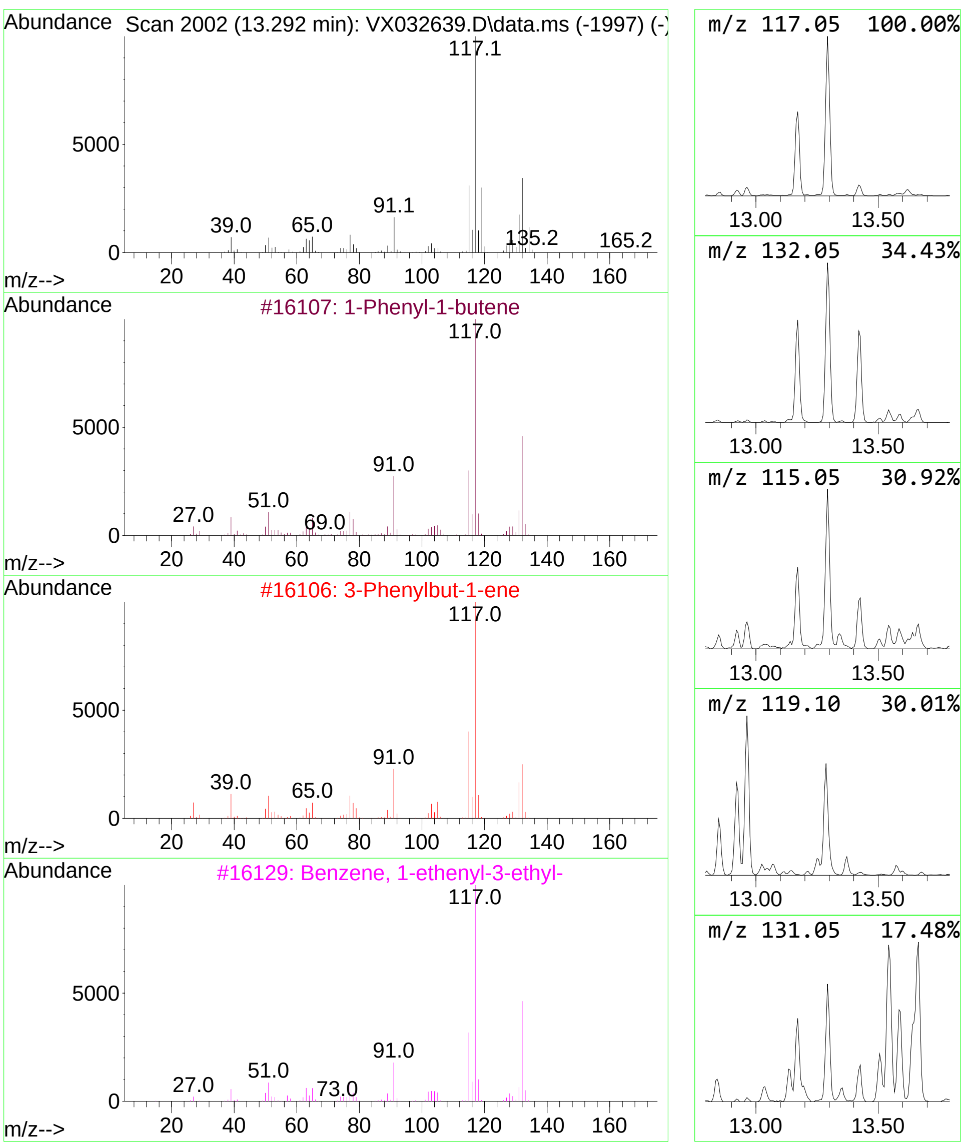


Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

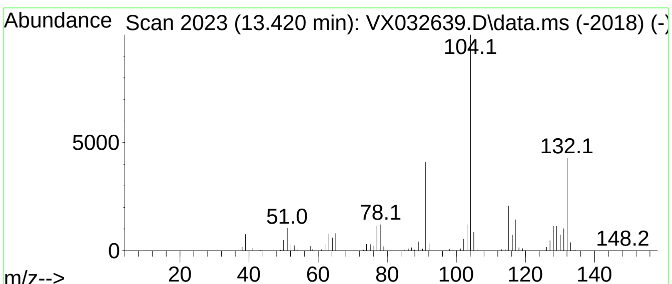
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

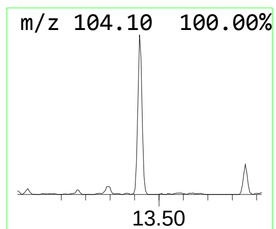
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

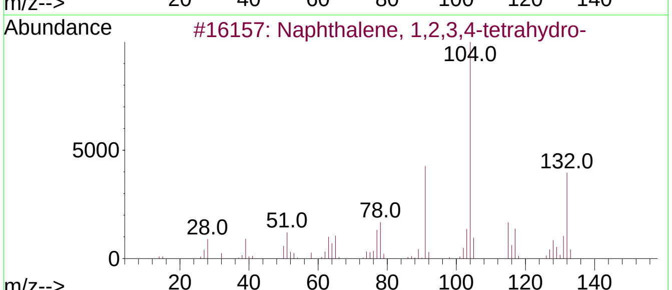
****
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

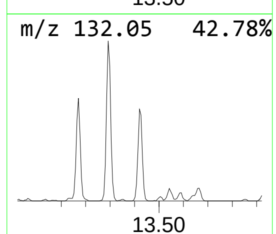
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	29.31 ug/l	133849	1,4-Dichlorobenzene-d4	12.024

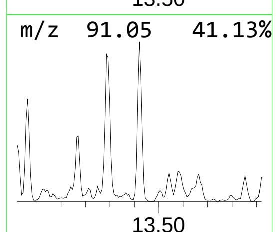
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Indan, epoxide	132	C9H8O	000768-22-9	49
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	49
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5	Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	35

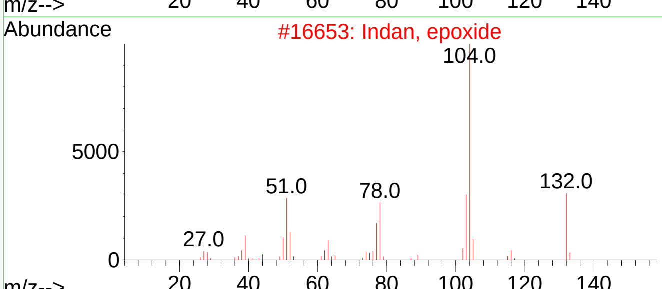


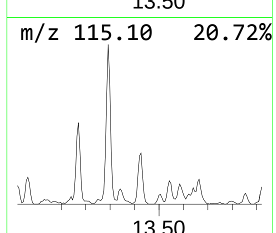


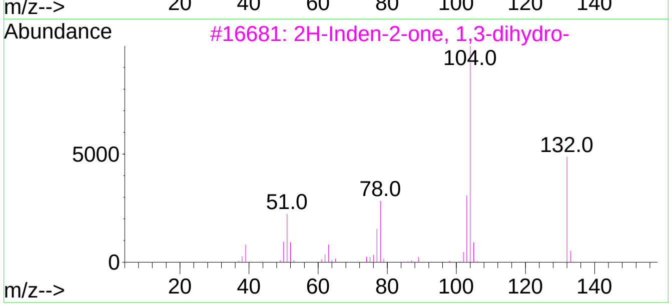


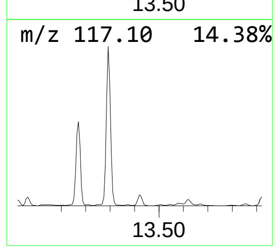












Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110722\
Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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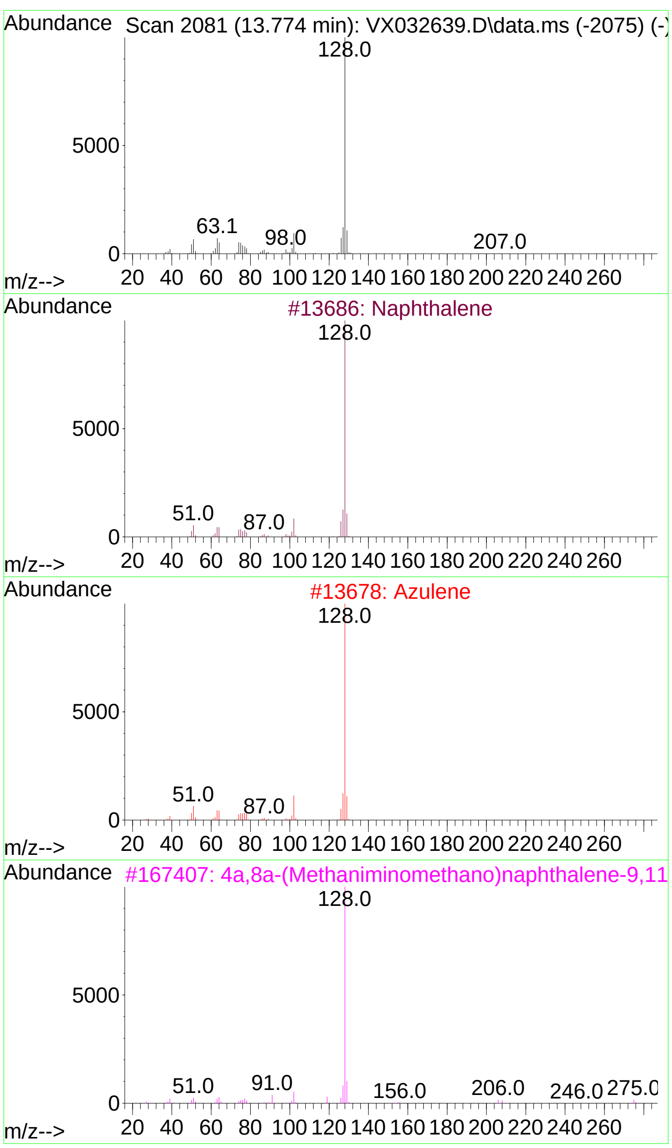
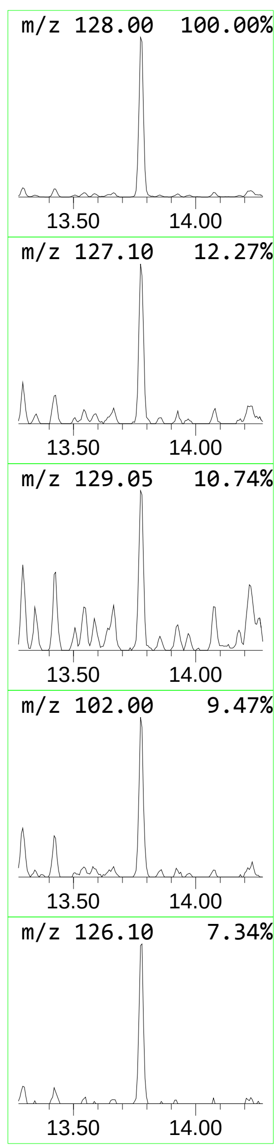
Peak Number 8 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.774	34.38 ug/l	157034	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	91
3	4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	64
4	1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	59
5	[4.2.2]Propella-2,4,7,9-tetraene	128	C10H8	088090-34-0	50





Library Search Compound Report

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Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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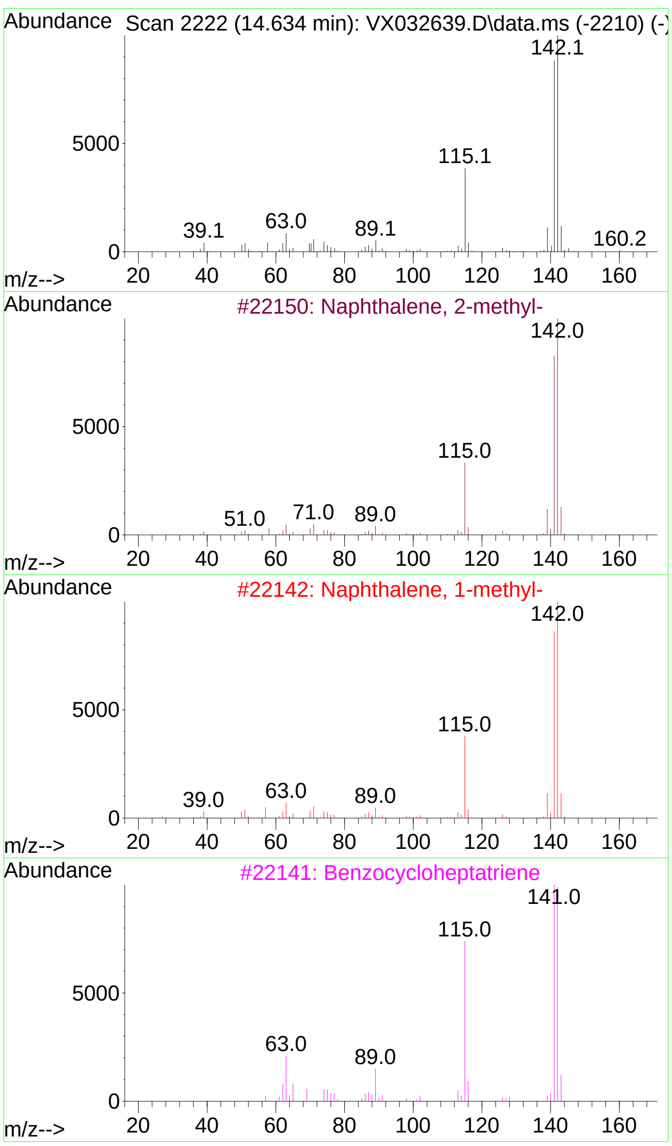
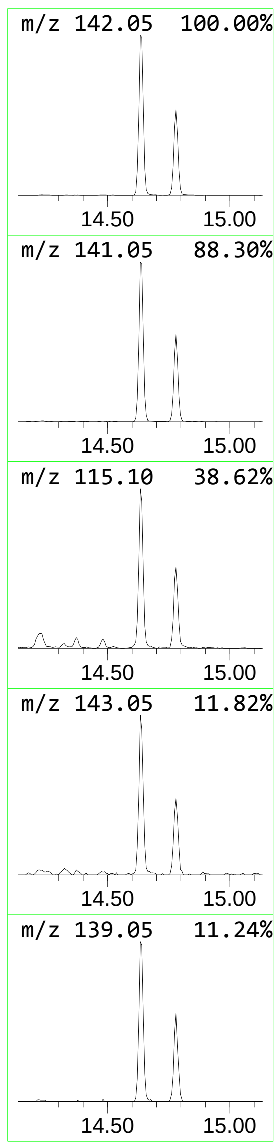
Peak Number 9 Benzocycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	77.84 ug/l	355536	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Library Search Compound Report

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Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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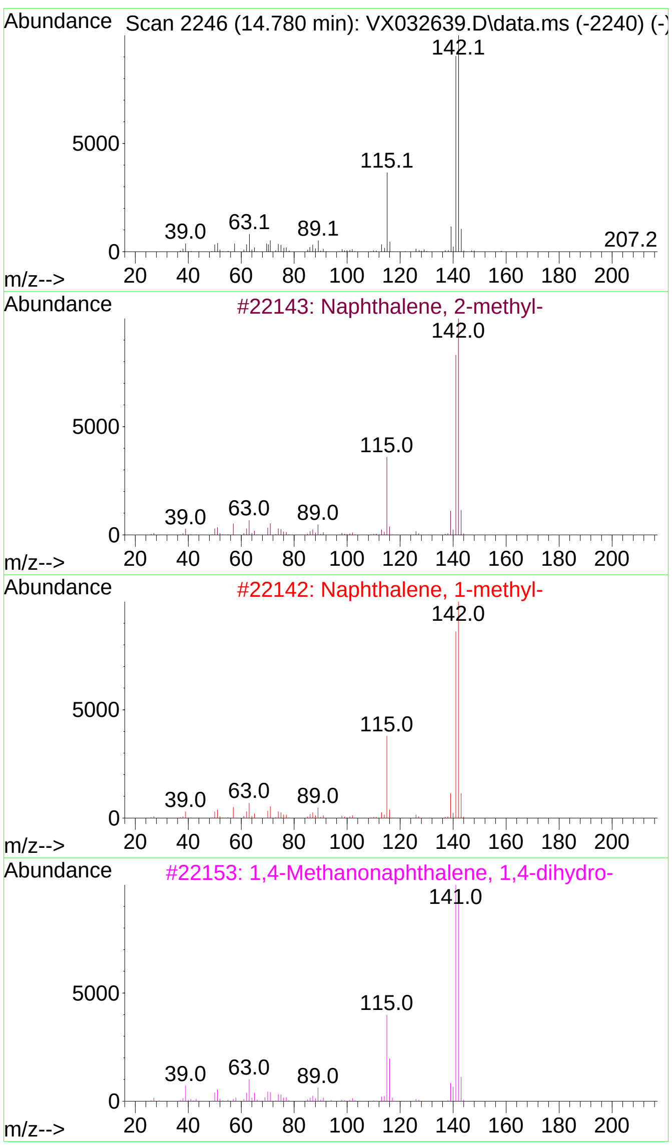
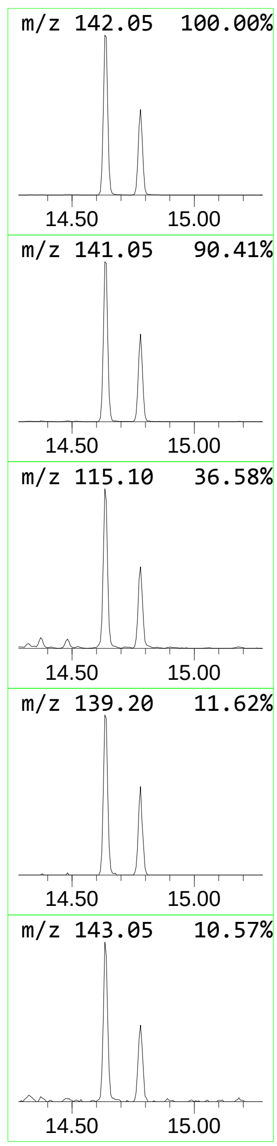
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	39.77 ug/l	181632	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86





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Data File : VX032639.D
Acq On : 07 Nov 2022 19:51
Operator : JC/MD
Sample : N5504-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SV34211L

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
Benzene, 1-ethy...	11.378	60.8	ug/l	277850	4	12.024	228366	50.0
Benzene, 1,2,3-...	12.085	43.3	ug/l	197921	4	12.024	228366	50.0
Indane	12.244	40.4	ug/l	184579	4	12.024	228366	50.0
Benzene, 1,2,3,...	12.963	27.5	ug/l	125515	4	12.024	228366	50.0
Benzene, 1-ethe...	13.171	28.1	ug/l	128094	4	12.024	228366	50.0
1-Phenyl-1-butene	13.292	64.8	ug/l	295781	4	12.024	228366	50.0
Naphthalene, 1,...	13.420	29.3	ug/l	133849	4	12.024	228366	50.0
Azulene	13.774	34.4	ug/l	157034	4	12.024	228366	50.0
Benzocyclohepta...	14.634	77.8	ug/l	355536	4	12.024	228366	50.0
1,4-Methanonaph...	14.780	39.8	ug/l	181632	4	12.024	228366	50.0