LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.386	764	774	790	rVB2	223581	508715	28.65%	1.925%
2	7.544	790	801	810	rBV	140043	263832	14.86%	0.998%
3	7.762	831	838	851	rBV2	206871	400684	22.56%	1.516%
4	8.232	911	918	927	rVB	251500	454717	25.61%	1.720%
5	8.349	932	938	943	rBV7	11986	23719	1.34%	0.090%
6	8.866	1016	1026	1030	rBV3	22296	43830	2.47%	0.166%
7	8.937	1032	1038	1051	rVB	264223	490339	27.61%	1.855%
8	9.172	1069	1078	1084	rBV9	9047	24836	1.40%	0.094%
9	9.331	1100	1105	1112	rVV7	12833	28754	1.62%	0.109%
10	9.413	1112	1119	1130	rVV2	228914	446606	25.15%	1.690%
11	9.583	1141	1148	1154	rVB9	10053	25737	1.45%	0.097%
12	9.707	1166	1169	1176	rVB8	14998	35604	2.01%	0.135%
13	9.783	1176	1182	1192	rBV4	24409	52145	2.94%	0.197%
14	10.136	1228	1242	1252	rBV2	202127	420214	23.66%	1.590%
15	10.218	1252	1256	1265	rBV2	14301	45667	2.57%	0.173%
16	10.288	1265	1268	1279	rVB9	13859	35999	2.03%	0.136%
17	10.470	1288	1299	1307	rBV5	39436	118137	6.65%	0.447%
18	10.611	1319	1323	1327	rVB4	17541	27942	1.57%	0.106%
19	10.682	1328	1335	1351	rVV2	294535	574837	32.37%	2.175%
20	10.987	1379	1387	1389	rBV7	19257	39335	2.22%	0.149%
21	11.058	1393	1399	1407	rBV	354796	632137	35.60%	2.392%
22	11.158	1412	1416	1420	rBV5	19711	31918	1.80%	0.121%
23	11.234	1420	1429	1433	rBV5	17461	53779	3.03%	0.203%
24	11.287	1434	1438	1455	rVB4	66497	219618	12.37%	0.831%
25	11.422	1455	1461	1463	rBV3	80633	139521	7.86%	0.528%
26	11.458	1463	1467	1474	rVV3	151659	311375	17.54%	1.178%
27	11.528	1474	1479	1486	rVB4	59476	104688	5.90%	0.396%
28	11.640	1493	1498	1504	rVB10	17080	26163	1.47%	0.099%
29	11.716	1506	1511	1521	rBV9	24678	52767	2.97%	0.200%
30	11.816	1521	1528	1532	rBV7	23327	42346	2.38%	0.160%
31	11.875	1534	1538	1542	rVB5	18481	27123	1.53%	0.103%
32	11.957	1544	1552	1556	rBV7	21222	56997	3.21%	0.216%
33	12.004	1556	1560	1567	rVB5	47699	85254	4.80%	0.323%
34	12.086	1567	1574	1579	rBV4	43819	102025	5.75%	0.386%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Title : SVOA CALIBRATION

35	12.145	1579	1584	1587	rVV7	23306	50061	2.82%	0.189%
36	12.186	1587	1591	1594	rVV4	35833	52008	2.93%	0.197%
37	12.257	1594	1603	1610	rVB5	59558	159920	9.01%	0.605%
38	12.356	1610	1620	1626	rBV8	38394	91529	5.15%	0.346%
39	12.415	1626	1630	1635	rBV6	26750	58683	3.30%	0.222%
40	12.509	1642	1646	1651	rBV2	35867	51450	2.90%	0.195%
41	12.568	1652	1656	1659	rBV4	49955	79310	4.47%	0.300%
42	12.697	1674	1678	1680	rBV	57388	76831	4.33%	0.291%
43	12.738	1680	1685	1688	rVV2	103963	147666	8.32%	0.559%
44	12.774	1689	1691	1696	rVB5	45309	61098	3.44%	0.231%
45	12.838	1696	1702	1703	rBV4	51933	88887	5.01%	0.336%
46	12.915	1711	1715	1722	rVB	113800	183917	10.36%	0.696%
47	13.003	1723	1730	1739	rBV2	92331	190187	10.71%	0.720%
48	13.091	1739	1745	1753	rBV7	60500	123853	6.97%	0.469%
49	13.208	1760	1765	1770	rVB8	36470	65112	3.67%	0.246%
50	13.267	1770	1775	1778	rBV7	45258	74890	4.22%	0.283%
51	13.344	1784	1788	1791	rBV3	60043	96646	5.44%	0.366%
52	13.443	1802	1805	1816	rVB9	31853	87531	4.93%	0.331%
53	13.579	1817	1828	1832	rVV5	63399	147779	8.32%	0.559%
54	13.626	1832	1836	1843	rVV5	45347	111753	6.29%	0.423%
55	13.690	1843	1847	1850	rVB3	20324	32112	1.81%	0.121%
56	13.731	1851	1854	1859	rVB6	29456	51086	2.88%	0.193%
57	13.802	1859	1866	1869	rBV4	62405	134497	7.57%	0.509%
58	13.896	1877	1882	1888	rVB7	56932	148325	8.35%	0.561%
59	14.037	1896	1906	1918	rVB6	103157	328984	18.53%	1.245%
60	14.178	1918	1930	1935	rBV3	112590	286942	16.16%	1.086%
61	14.248	1935	1942	1946	rBV	545079	919292	51.77%	3.478%
62	14.289	1947	1949	1954	rVB2	101795	136060	7.66%	0.515%
63	14.413	1967	1970	1974	rVB4	34391	45630	2.57%	0.173%
64	14.489	1980	1983	1987	rVB3	62770	83566	4.71%	0.316%
65	14.554	1987	1994	2007	rBV	578756	1061689	59.79%	4.017%
66	14.718	2017	2022	2025	rBV7	29889	40644	2.29%	0.154%
67	14.765	2026	2030	2033	rBV6	24308	26964	1.52%	0.102%
68	14.812	2033	2038	2040	rBV2	142661	209860	11.82%	0.794%
69	14.854	2040	2045	2055	rVB	610728	1061868	59.80%	4.017%
70	15.065	2073	2081	2091	rBV3	227600	551400	31.05%	2.086%
71	15.147	2091	2095	2100	rVB2	89251	142808	8.04%	0.540%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Title : SVOA CALIBRATION

72	15.235	2105	2110	2118	rVB10	46122	126563	7.13%	0.479%
73	15.312	2118	2123	2127	rBV4	32347	68206	3.84%	0.258%
74	15.359	2128	2131	2142	rVB9	47553	88580	4.99%	0.335%
75	15.459	2144	2148	2152	rBV5	33250	56791	3.20%	0.215%
76	15.512	2153	2157	2166	rVB8	70494	154492	8.70%	0.585%
77	15.611	2168	2174	2191	rVB3	102313	277499	15.63%	1.050%
78	15.841	2201	2213	2228	rBV	759147	1374733	77.42%	5.201%
79	15.970	2228	2235	2241	rBV2	267045	437707	24.65%	1.656%
80	16.046	2246	2248	2252	rVB4	35344	40680	2.29%	0.154%
81	16.428	2309	2313	2322	rVB7	34133	74890	4.22%	0.283%
82	16.505	2322	2326	2329	rBV6	45328	79603	4.48%	0.301%
83	16.546	2330	2333	2338	rVB	58571	85116	4.79%	0.322%
84	16.728	2357	2364	2368	rVB10	25521	42081	2.37%	0.159%
85	16.810	2373	2378	2384	rBV2	124695	170295	9.59%	0.644%
86	16.875	2386	2389	2391	rBV4	21985	29814	1.68%	0.113%
87	16.951	2398	2402	2407	rVB8	16820	24387	1.37%	0.092%
88	17.298	2453	2461	2465	rVB8	31687	69834	3.93%	0.264%
89	17.439	2481	2485	2490	rVB4	20802	30683	1.73%	0.116%
90	17.597	2505	2512	2522	rBV2	842677	1344661	75.72%	5.087%
91	17.697	2522	2529	2536	rVB2	842993	1322511	74.48%	5.004%
92	17.768	2537	2541	2547	rVB4	53538	87132	4.91%	0.330%
93	18.032	2580	2586	2596	rVB4	27972	72445	4.08%	0.274%
94	18.438	2650	2655	2660	rBV7	53182	78331	4.41%	0.296%
95	19.971	2909	2916	2928	rBV	1200404	1775731	100.00%	6.718%
96	21.892	3235	3243	3256	rVB	970848	1755767	98.88%	6.643%
97	25.077	3774	3785	3796	rBV3	508932	1574983	88.69%	5.959%
98	25.306	3814	3824	3842	rVB2	474493	1524951	85.88%	5.770%
99	26.393	4002	4009	4020	rVB6	75352	220400	12.41%	0.834%
100	27.815	4243	4251	4265	rVB7	56541	234078	13.18%	0.886%

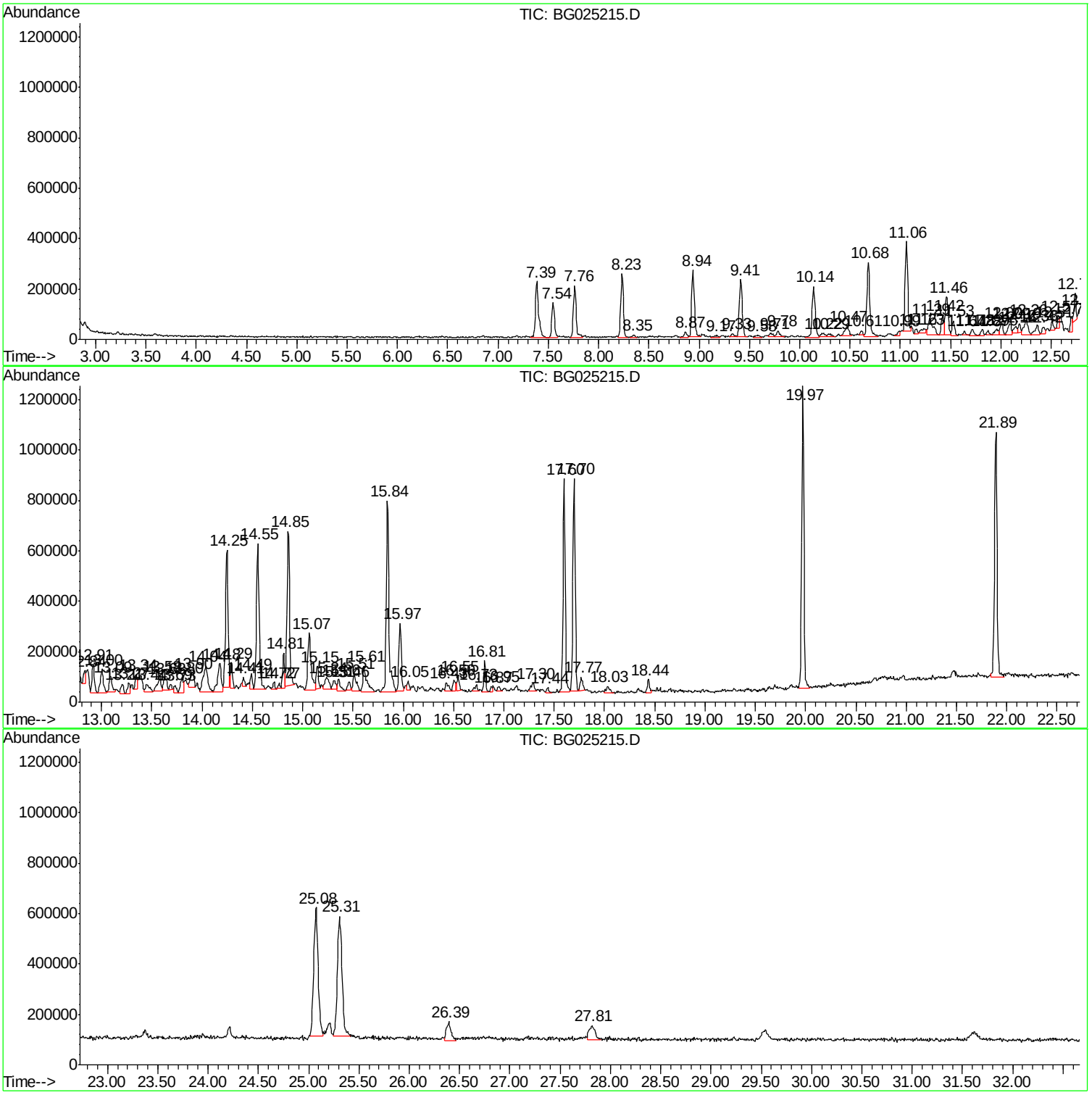
Sum of corrected areas: 26431142

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

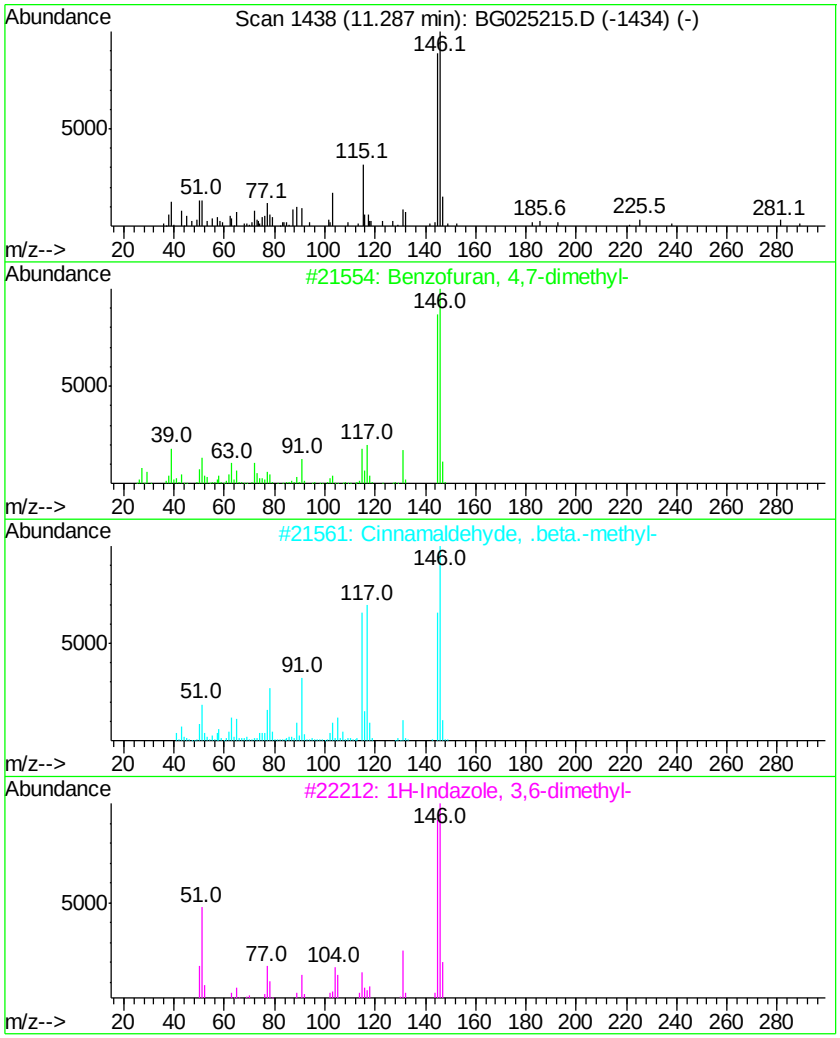
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

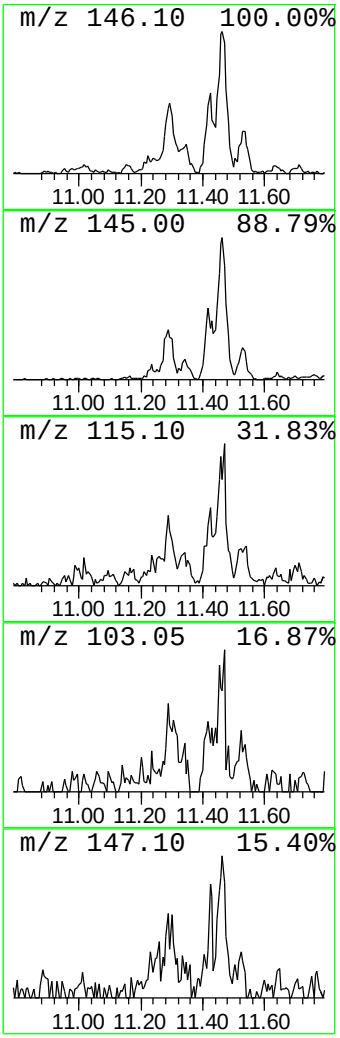
****

Peak Number 2 Benzofuran, 4,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	6.95 ng/ul	219618	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	74
2	Cinnamaldehyde, .beta.-methyl-	146	C10H10O	001196-67-4	64
3	1H-Indazole, 3,6-dimethyl-	146	C9H10N2	007746-28-3	64
4	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	64
5	Benzofuran-2-carboxaldehyde	146	C9H6O2	004265-16-1	59





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

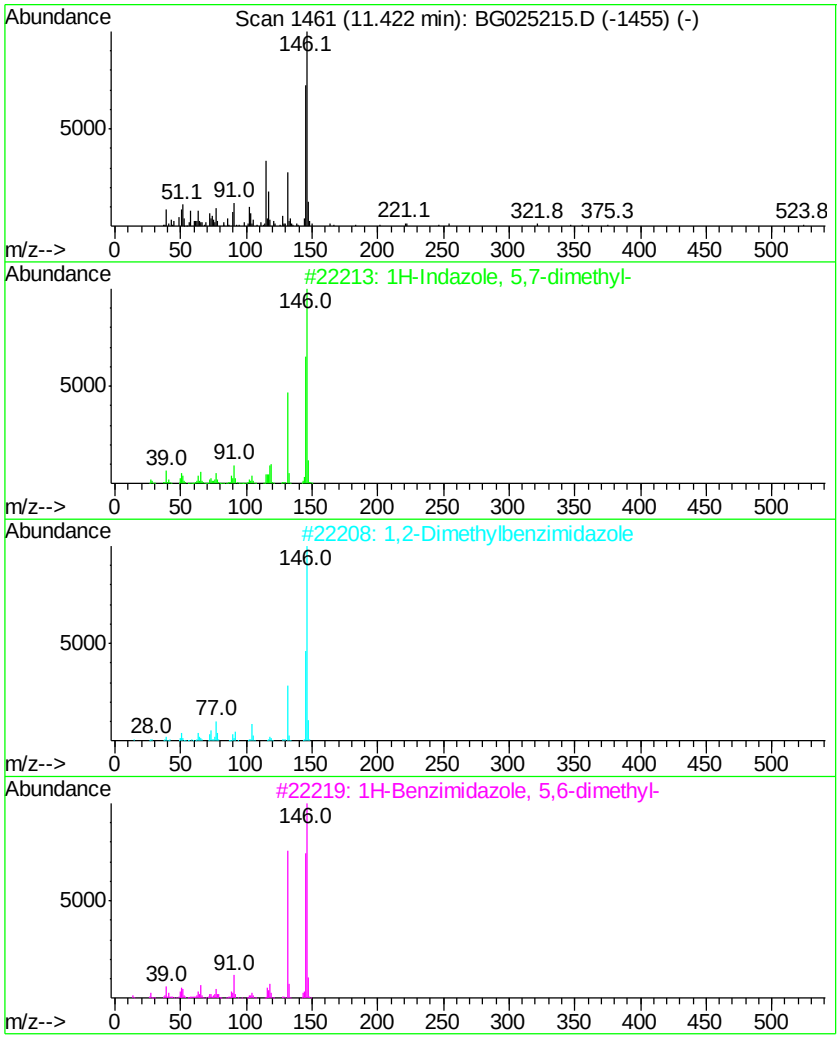
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

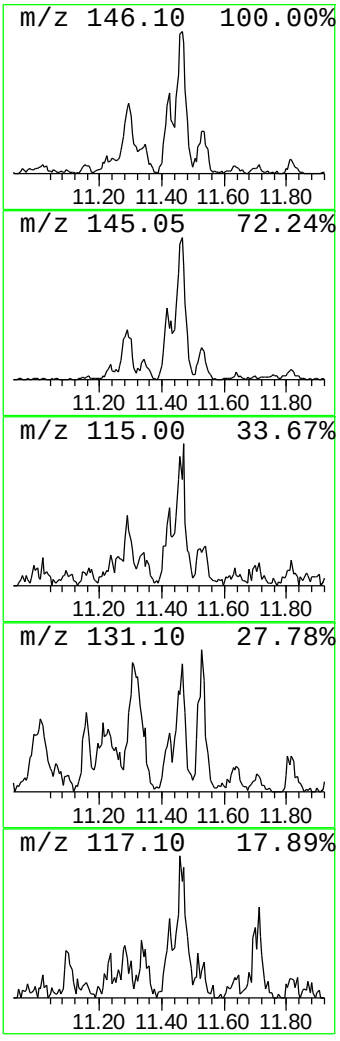
****

Peak Number 3 1H-Indazole, 5,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	4.41 ng/ul	139521	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	68
2	1,2-Dimethylbenzimidazole	146	C9H10N2	002876-08-6	64
3	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	64
4	Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	59
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	53





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

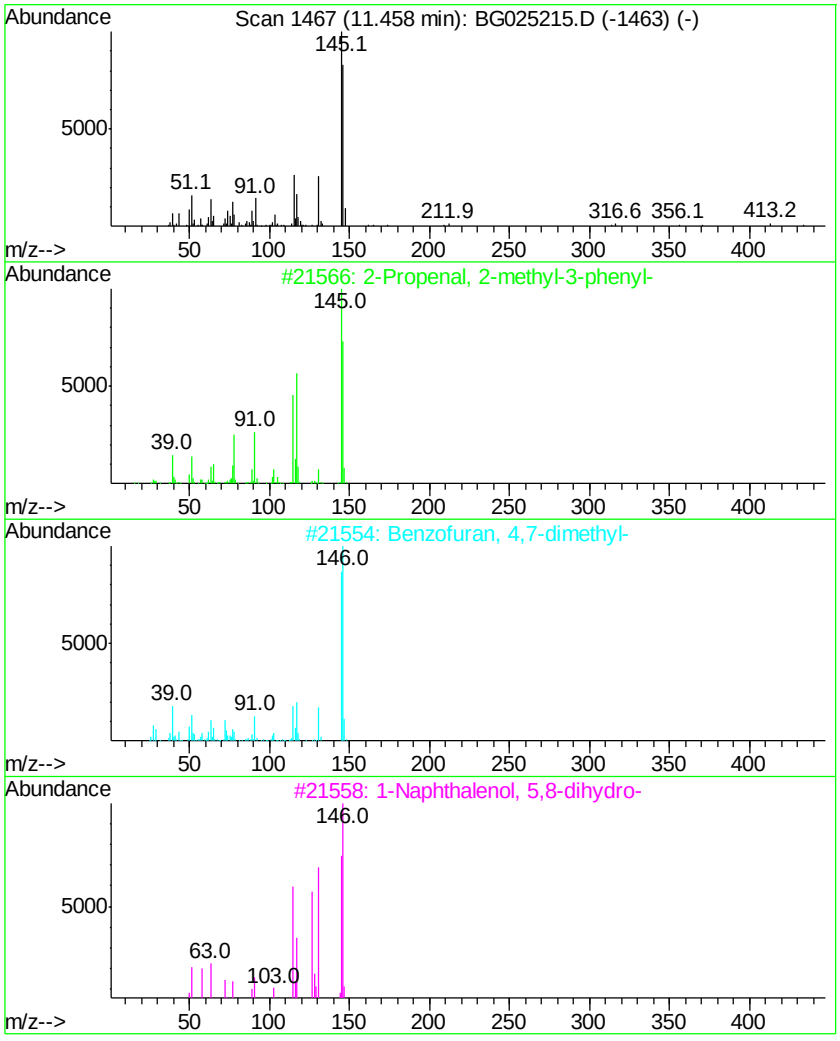
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

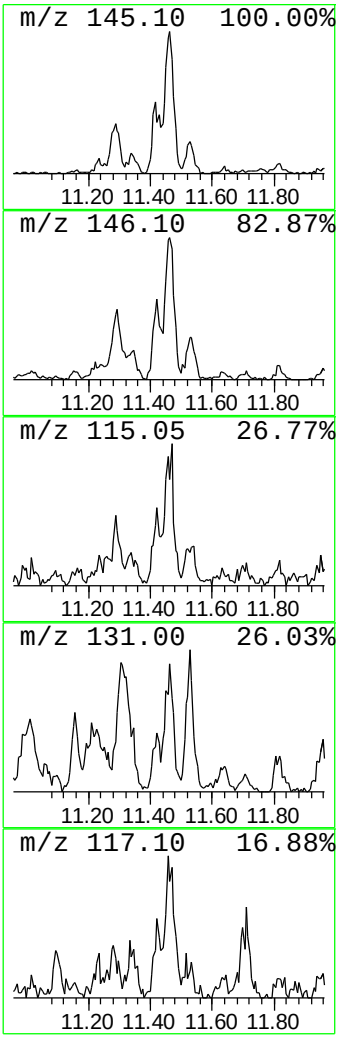
****

Peak Number 4 2-Propenal, 2-methyl-3-phenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	9.85 ng/ul	311375	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 2-methyl-3-phenyl-	146	C10H10O	000101-39-3	59
2	Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	58
3	1-Naphthalenol, 5,8-dihydro-	146	C10H10O	027673-48-9	53
4	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	50
5	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	50





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

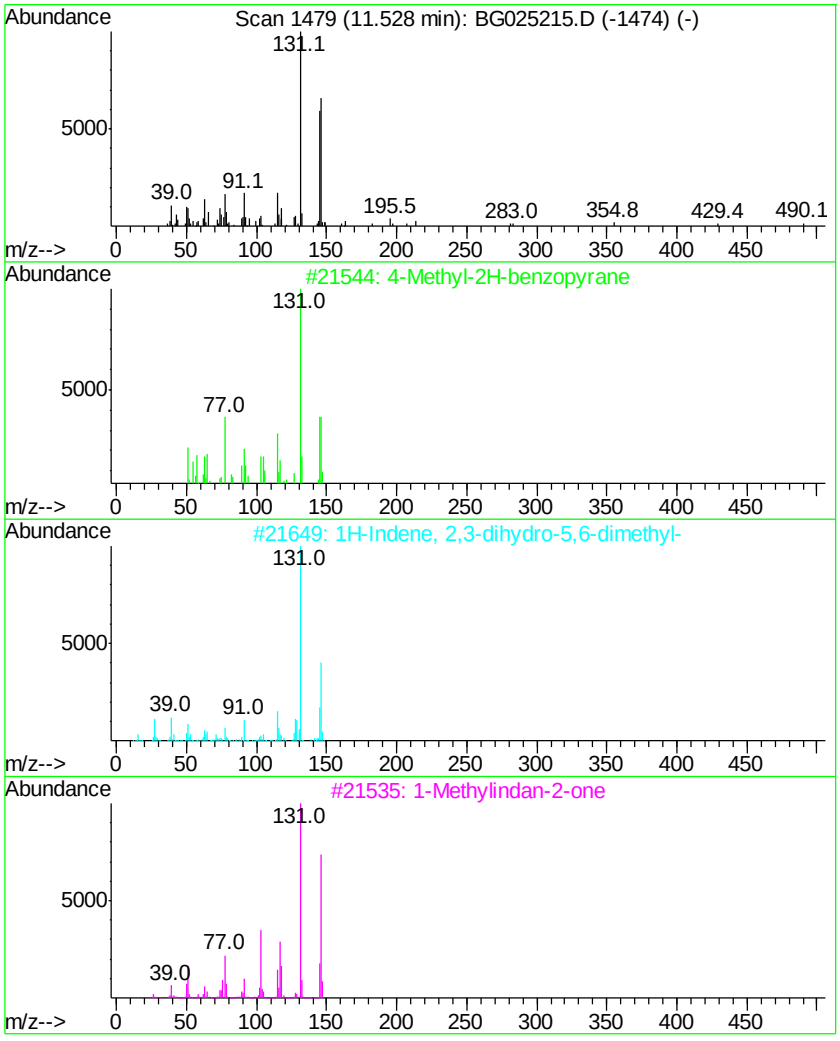
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

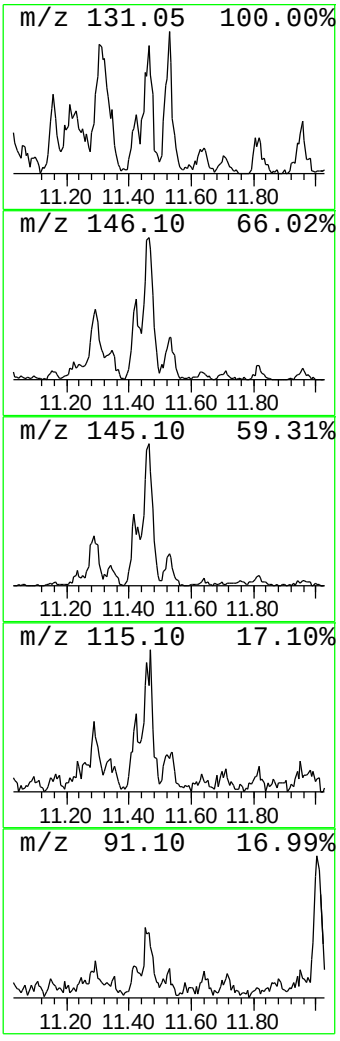
****

Peak Number 5 4-Methyl-2H-benzopyrane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	3.31 ng/ul	104688	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-2H-benzopyrane	146	C10H10O	021776-94-3	87
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	59
3	1-Methylindan-2-one	146	C10H10O	035587-60-1	53
4	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	53
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	52





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

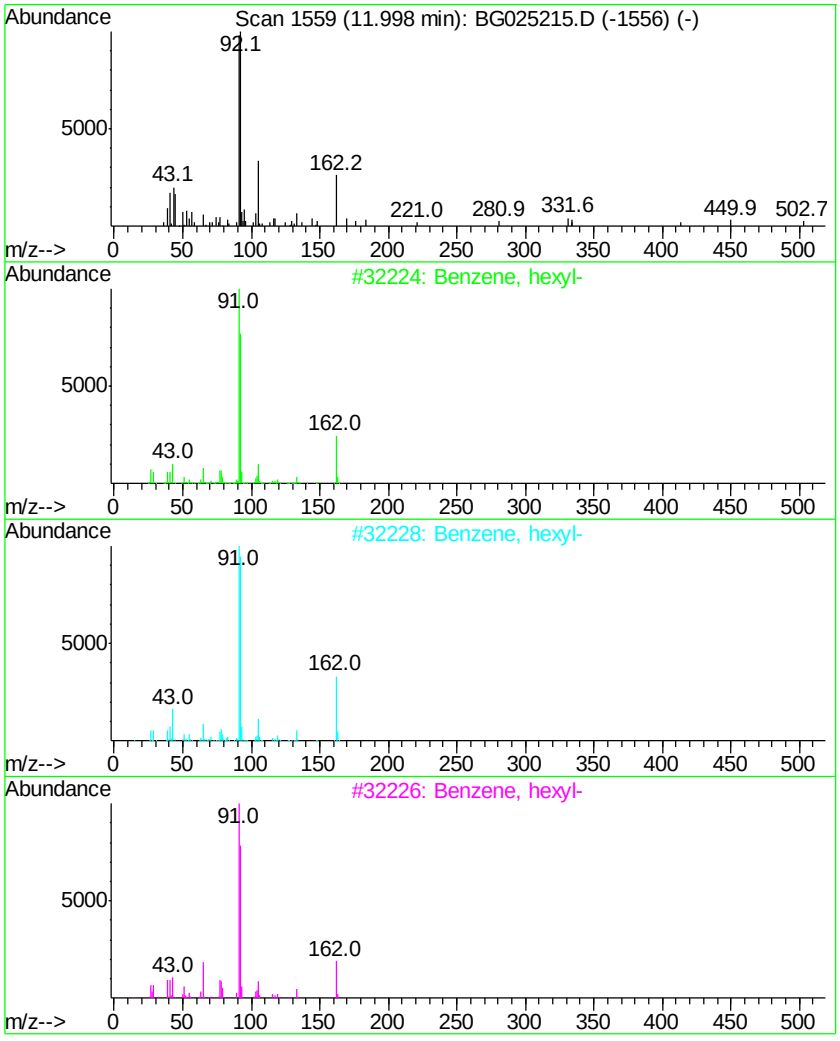
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

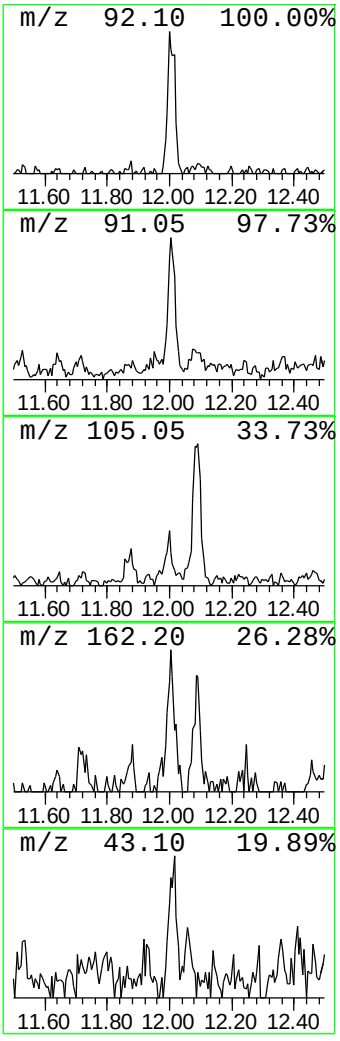
****

Peak Number 6 Benzene, hexyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.70 ng/ul	85254	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, hexyl-	162	C12H18	001077-16-3	83
2	Benzene, hexyl-	162	C12H18	001077-16-3	83
3	Benzene, hexyl-	162	C12H18	001077-16-3	80
4	8,9-Dimethylene-4-oxatricyclo[5....	162	C11H14O	080707-63-7	72
5	Bicyclo[6.4.0]dodeca-9,11-diene	162	C12H18	1000160-45-0	72





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

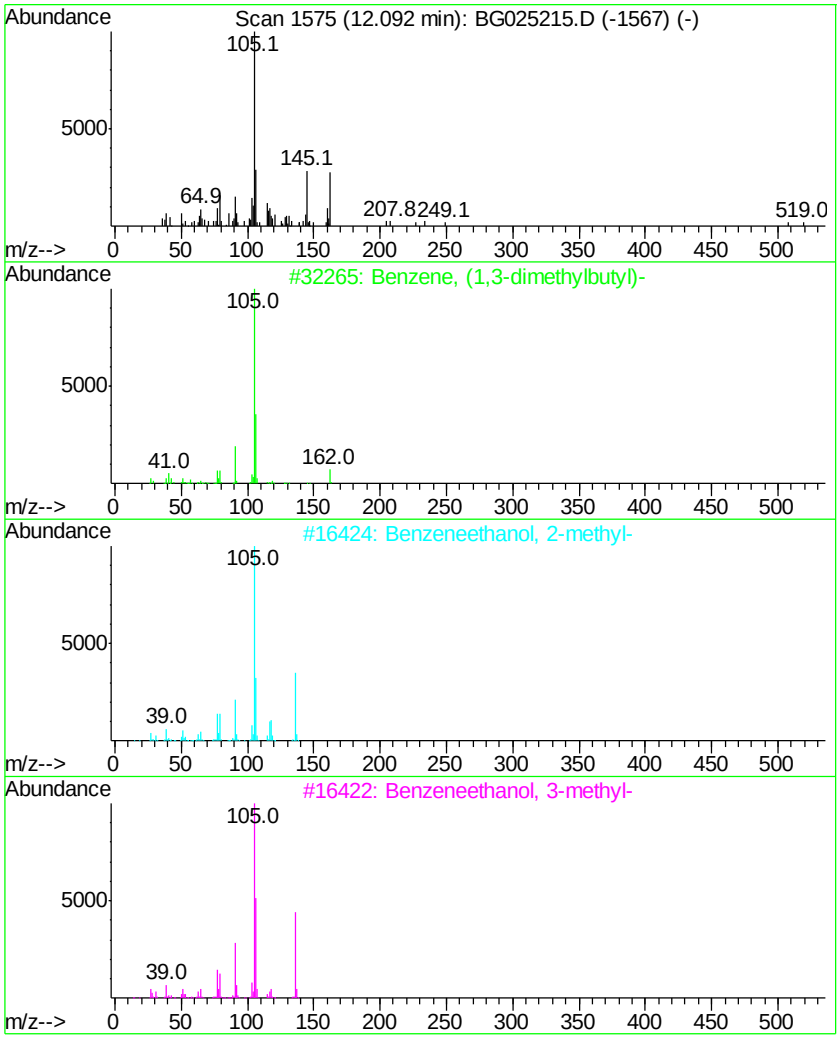
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

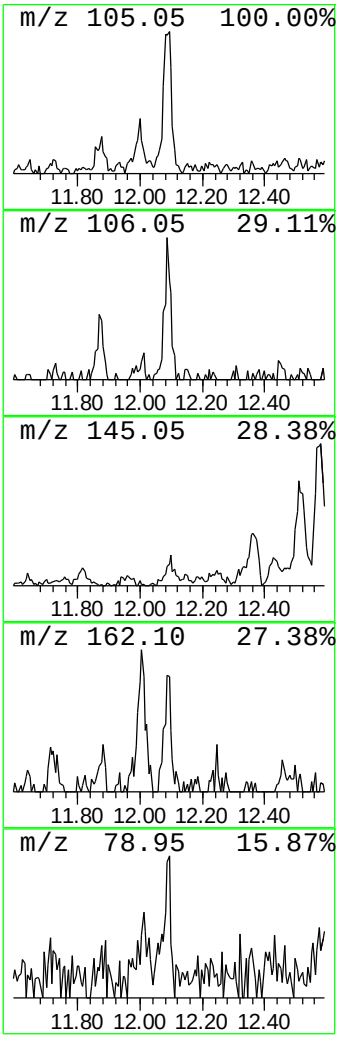
****

Peak Number 7 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	3.23 ng/ul	102025	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,3-dimethylbutyl)-	162	C12H18	019219-84-2	46
2	Benzeneethanol, 2-methyl-	136	C9H12O	019819-98-8	38
3	Benzeneethanol, 3-methyl-	136	C9H12O	001875-89-4	35
4	(1H)Pyrrole-3-carbonitrile, 2-me...	106	C6H6N2	026187-27-9	27
5	Benzene, (1,2-dimethyl-3-butenyl)-	160	C12H16	050871-04-0	22





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

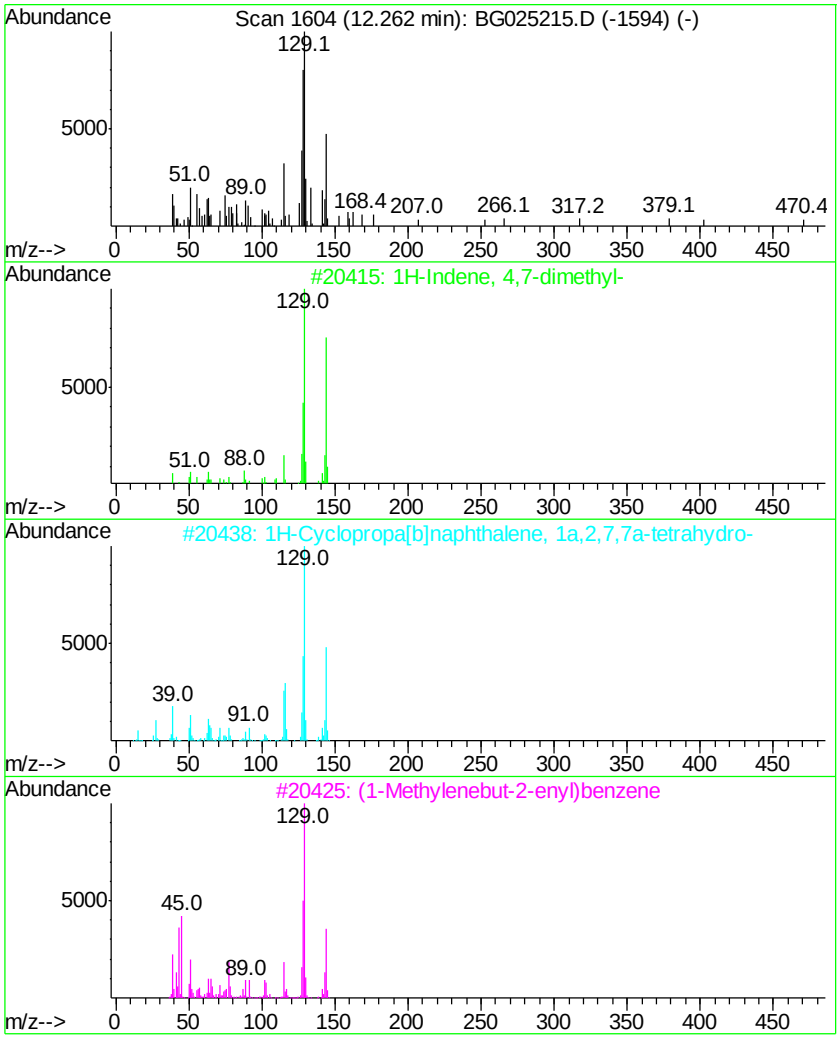
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

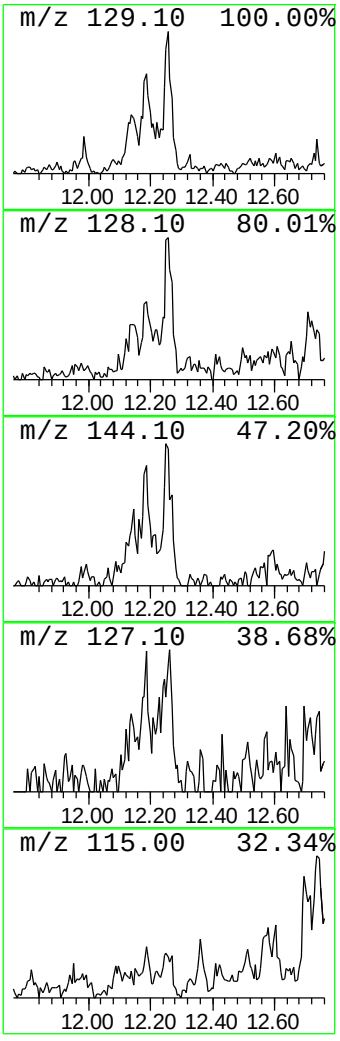
****

Peak Number 8 1H-Indene, 4,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	5.06 ng/ul	159920	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	87
2	1H-Cyclopropa[b]naphthalene, 1a,...	144	C11H12	006571-72-8	87
3	(1-Methylenebut-2-enyl)benzene	144	C11H12	070588-46-4	80
4	1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	80
5	Naphthalene, 1,2-dihydro-3-methyl-	144	C11H12	002717-44-4	76





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

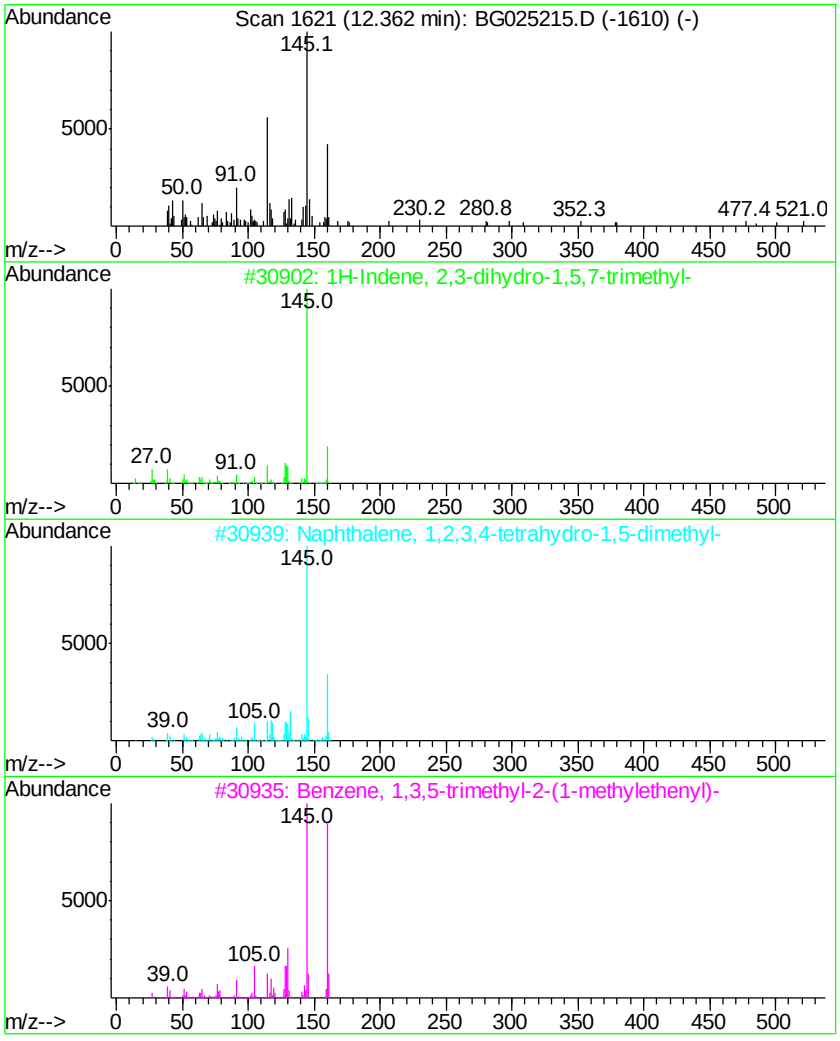
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

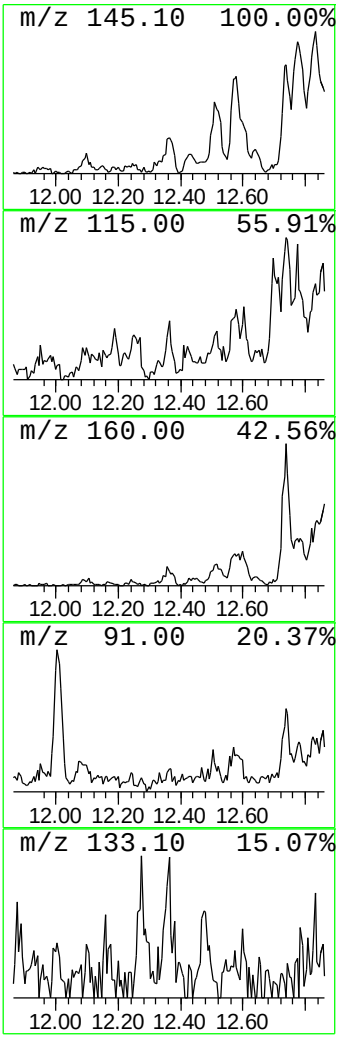
****

Peak Number 9 1H-Indene, 2,3-dihydro-1,5,... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	2.90 ng/ul	91529	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,5,7-tri...	160	C12H16	054340-88-4	59
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	50
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	49
4	Silane, methyltripropynyl-	160	C10H12Si	1000158-62-1	47
5	Ethanone, 1-(2-benzofuranyl)-	160	C10H8O2	001646-26-0	47





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

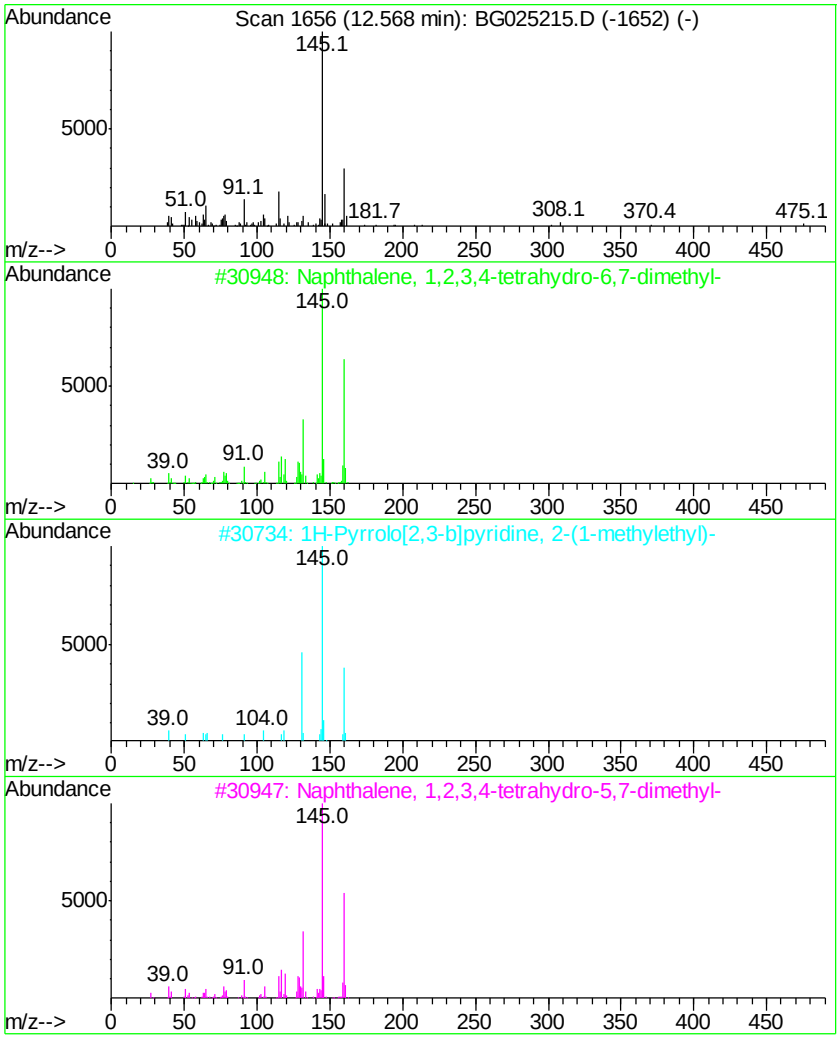
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

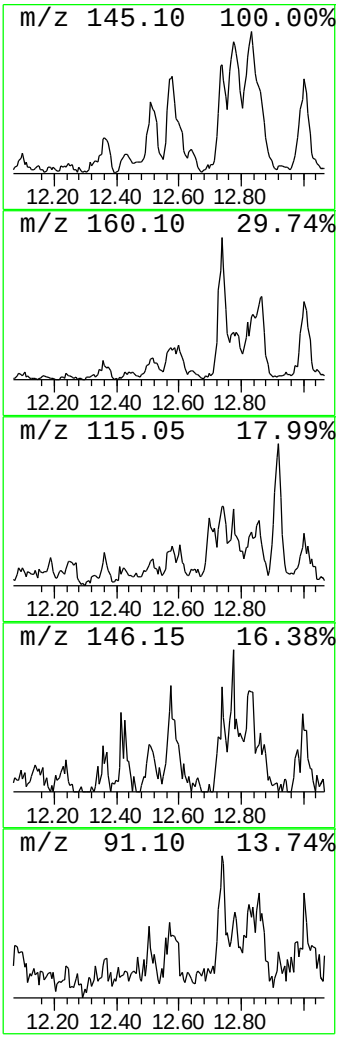
****

Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	2.51 ng/ul	79310	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	86
2	1H-Pyrrolo[2,3-b]pyridine, 2-(1-...	160	C10H12N2	027257-18-7	86
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	86
4	Ethanone, 1-(2,3-dihydro-1H-inde...	160	C11H12O	004228-10-8	80
5	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	80





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

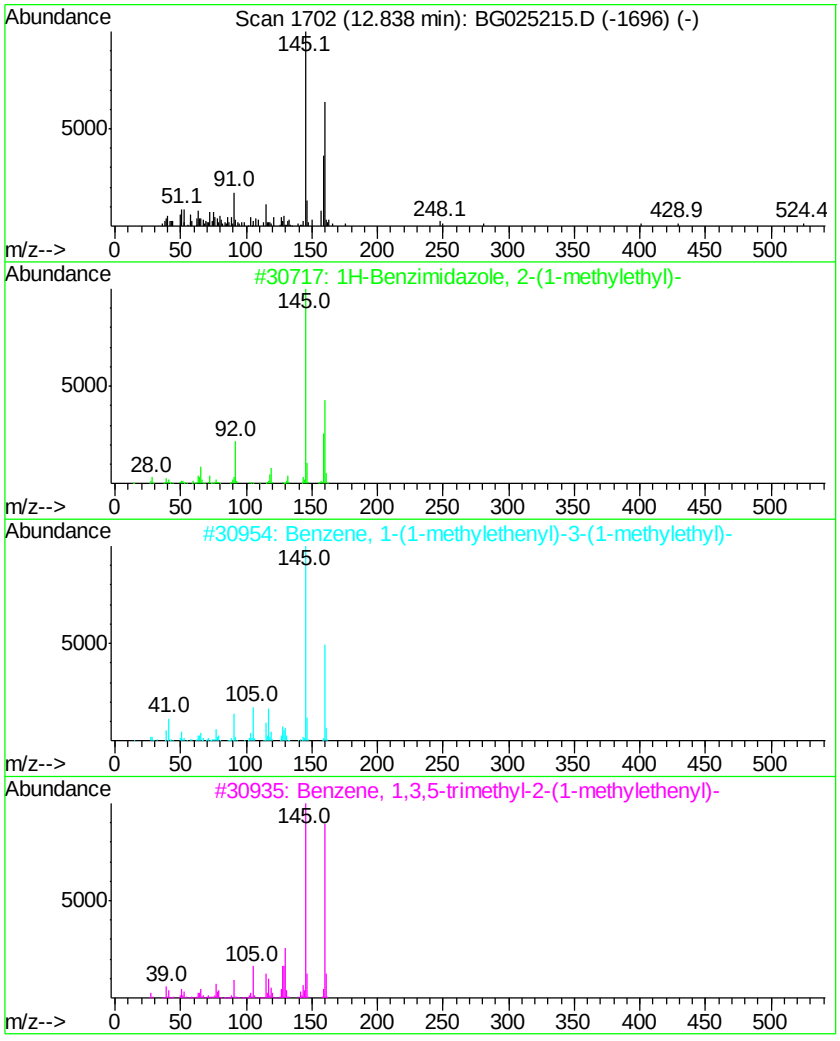
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

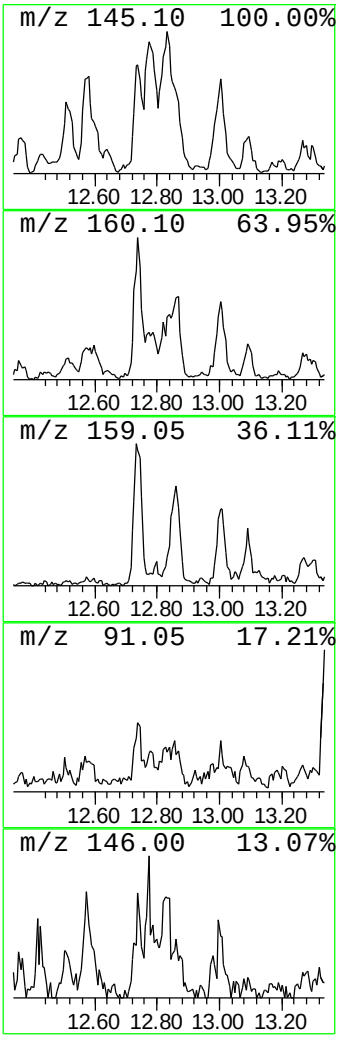
****

Peak Number 11 1H-Benzimidazole, 2-(1-meth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.84	2.81 ng/ul	88887	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-(1-methyleth...	160	C10H12N2	005851-43-4	86
2	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	72
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	64
4	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	64
5	Benzene, (2,2-dimethyl-1-methyle...	160	C12H16	005676-29-9	64





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

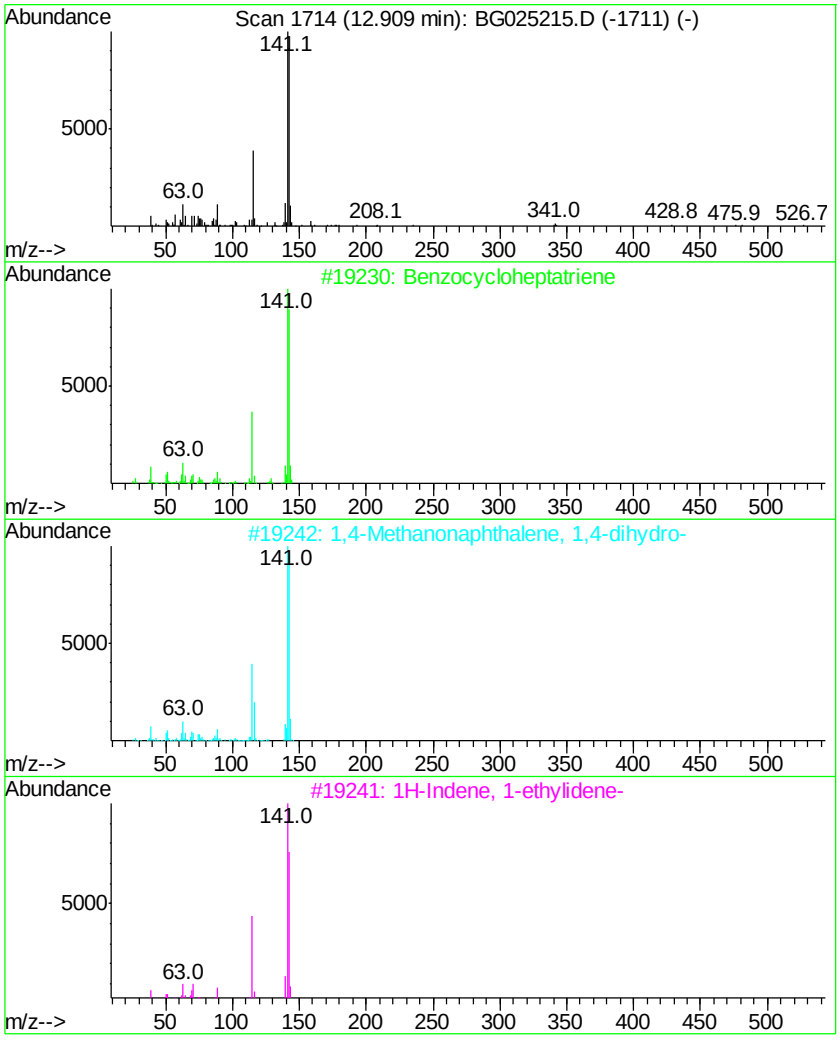
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

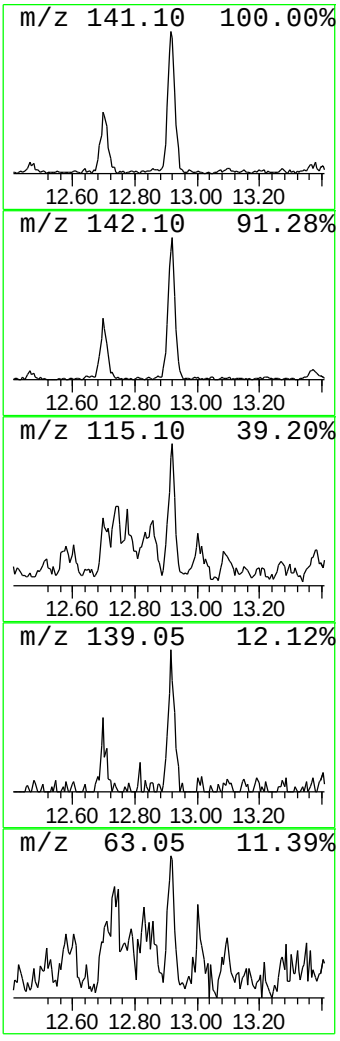
****

Peak Number 12 Benzocycloheptatriene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	5.82 ng/ul	183917	Naphthalene-d8	11.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzocycloheptatriene	142	C11H10	000264-09-5	94
2	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
3	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	94
4	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

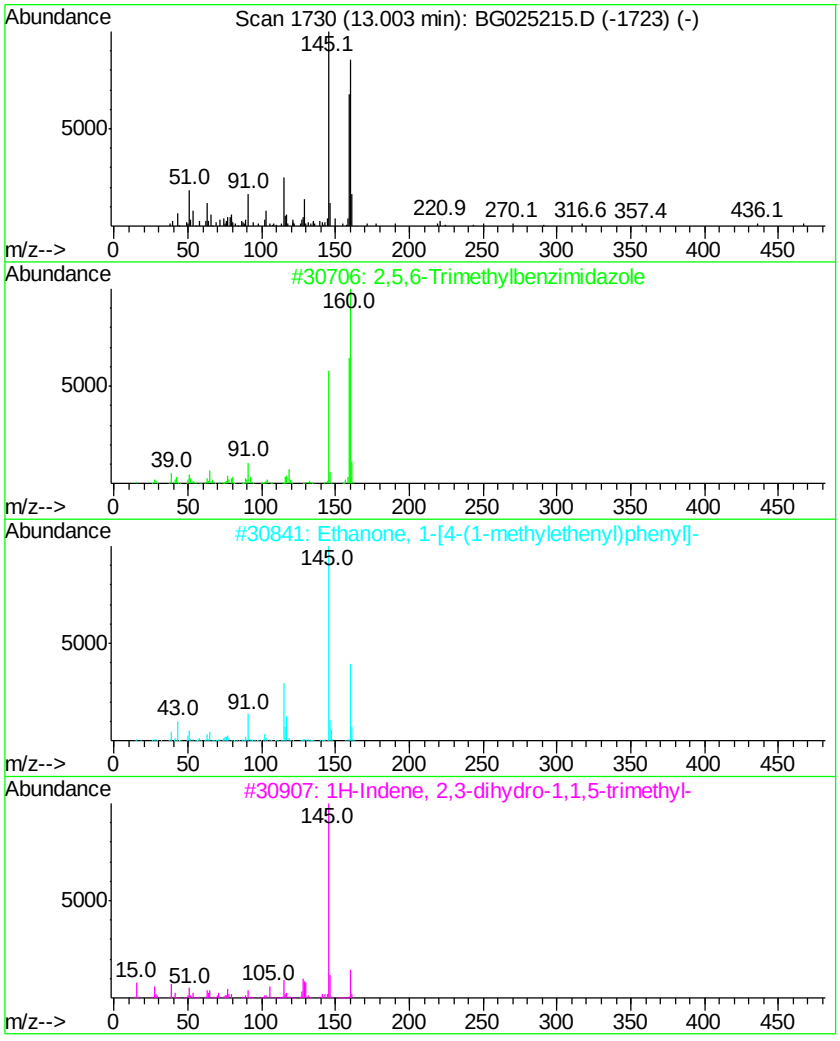
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

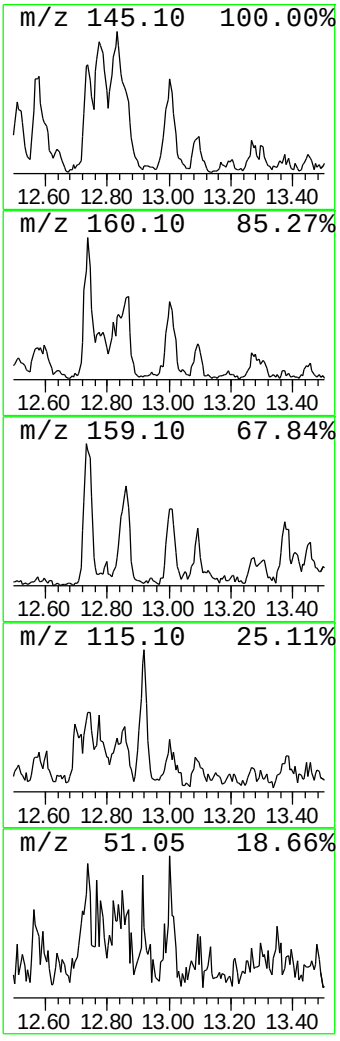
****

Peak Number 13 2,5,6-Trimethylbenzimidazole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	3.58 ng/ul	190187	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5,6-Trimethylbenzimidazole	160	C10H12N2	003363-56-2	80
2	Ethanone, 1-[4-(1-methylethenyl)...	160	C11H12O	005359-04-6	70
3	1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	64
4	Benzene, 1-(1-methylethenyl)-4-(...	160	C12H16	002388-14-9	62
5	1H-Indene, 2,3-dihydro-4,5,7-tri...	160	C12H16	006682-06-0	58





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

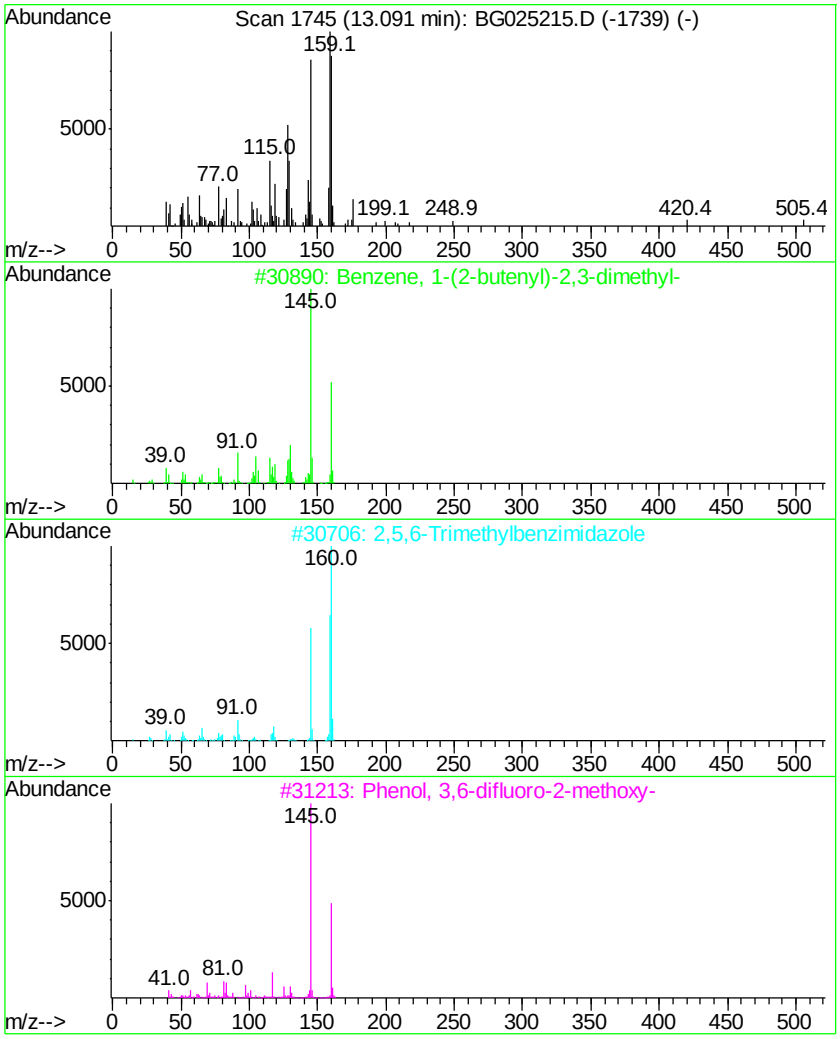
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

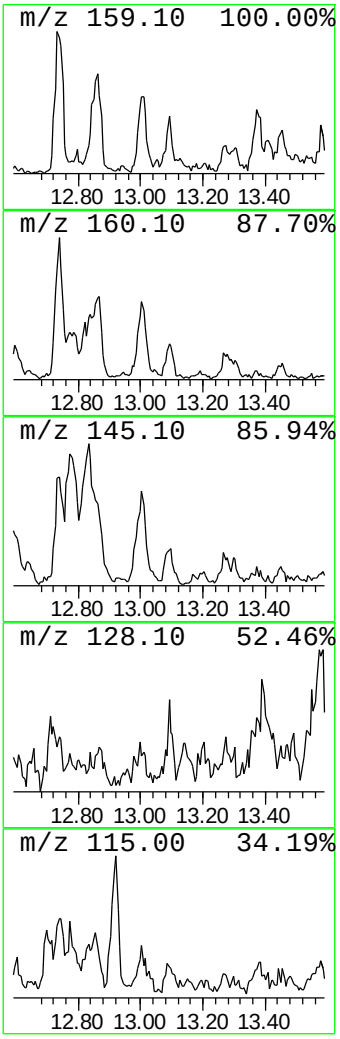
****

Peak Number 14 unknown-02 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	2.33 ng/ul	123853	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	43
2	2,5,6-Trimethylbenzimidazole	160	C10H12N2	003363-56-2	43
3	Phenol, 3,6-difluoro-2-methoxy-	160	C7H6F2O2	075626-22-1	38
4	2,3,4-Trifluorobenzaldehyde	160	C7H3F3O	161793-17-5	30
5	Oct-3-ene-1,5-diyne, 3-isopropyl...	160	C12H16	1000211-23-1	27





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

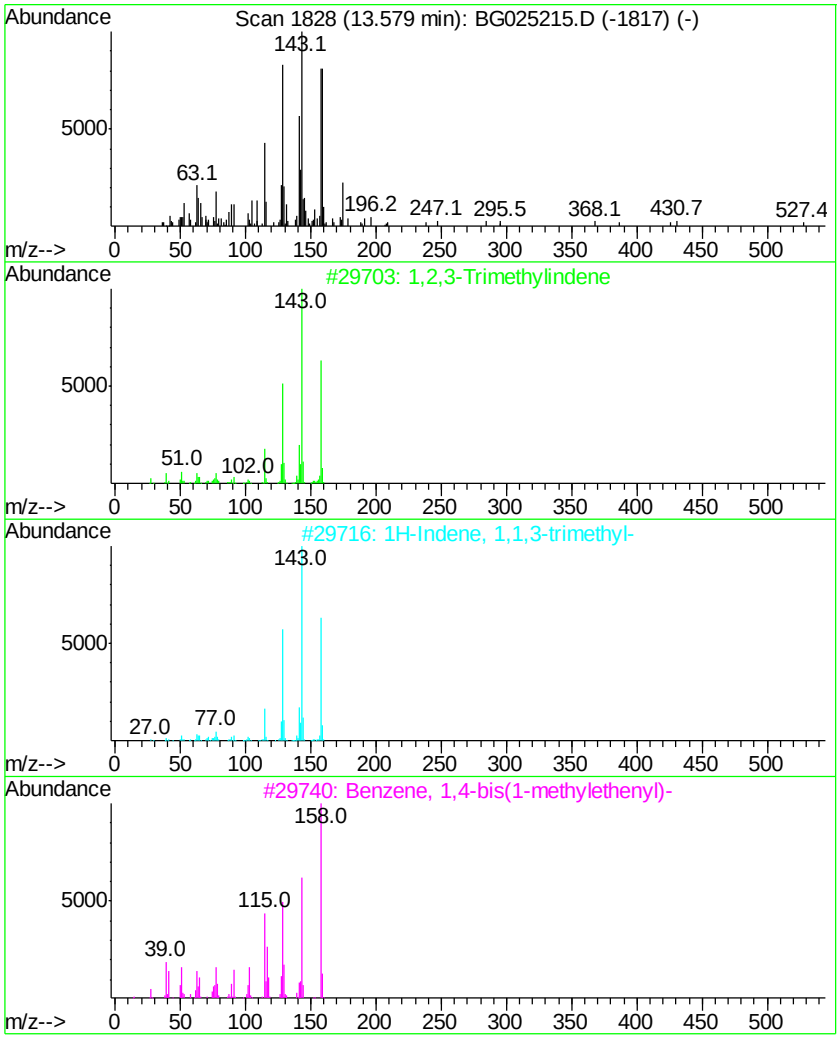
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

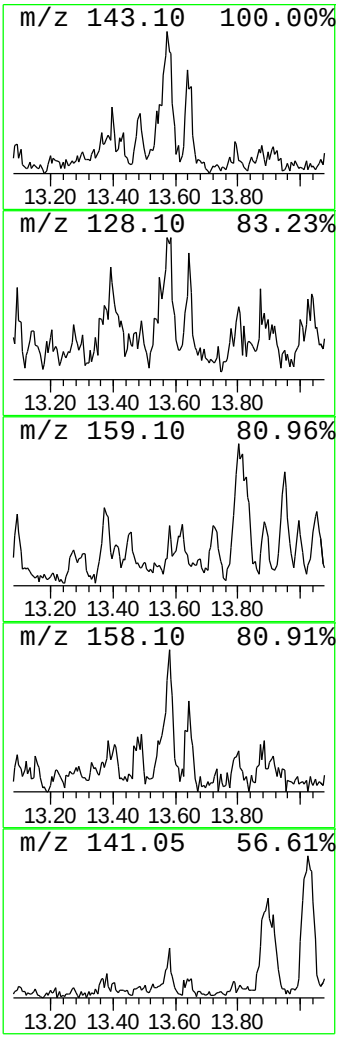
****

Peak Number 15 1,2,3-Trimethylindene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	2.78 ng/ul	147779	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3-Trimethylindene	158	C12H14	004773-83-5	70
2	1H-Indene, 1,1,3-trimethyl-	158	C12H14	002177-45-9	70
3	Benzene, 1,4-bis(1-methylethenyl)-	158	C12H14	001605-18-1	58
4	Benzene, 1,4-bis(1-methylethenyl)-	158	C12H14	001605-18-1	58
5	1,2,3-Trimethylindene	158	C12H14	004773-83-5	53





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

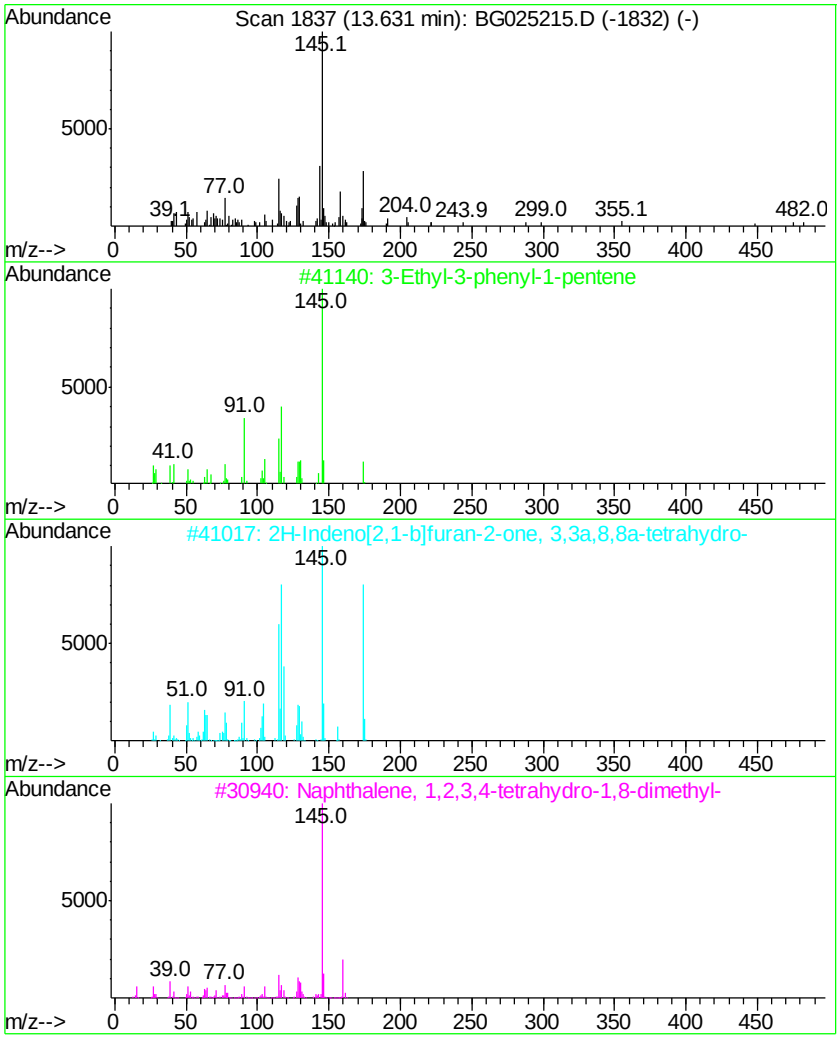
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

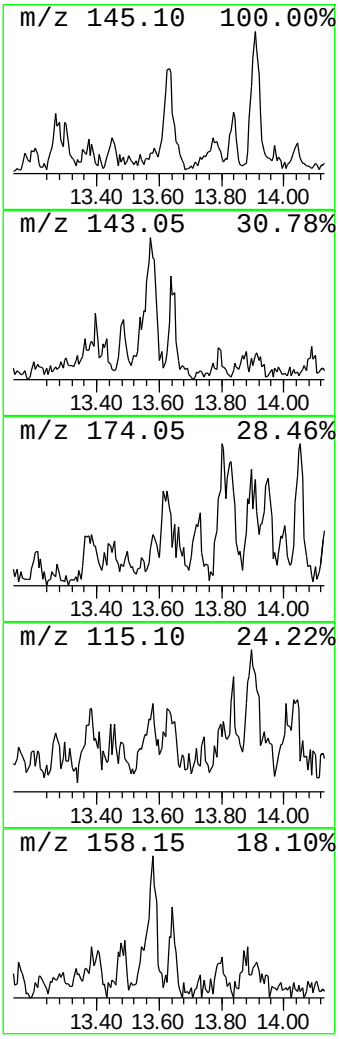
****

Peak Number 16 3-Ethyl-3-phenyl-1-pentene Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	2.10 ng/ul	111753	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Ethyl-3-phenyl-1-pentene	174	C13H18	019781-34-1	59
2	2H-Indeno[2,1-b]furan-2-one, 3,3...	174	C11H10O2	080343-98-2	47
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	47
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	042775-77-9	46
5	4-(N-Methyl-N-methoxy)indancarbo...	205	C12H15NO2	1000284-45-9	46





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

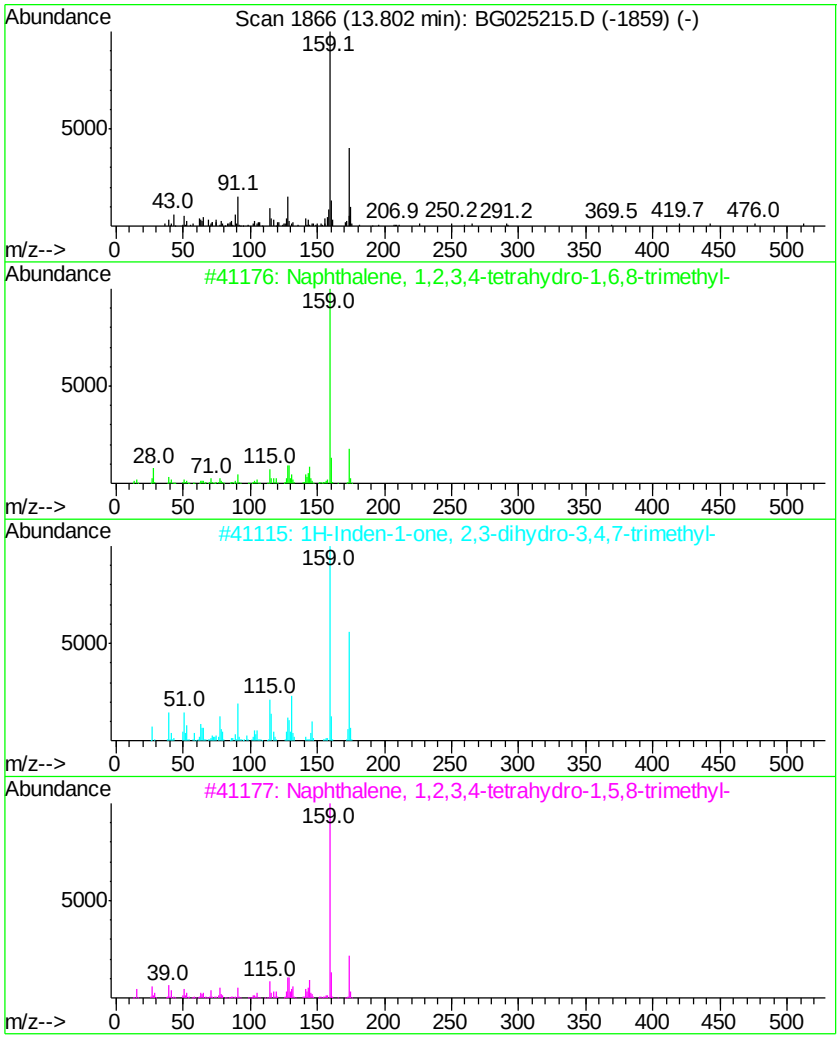
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

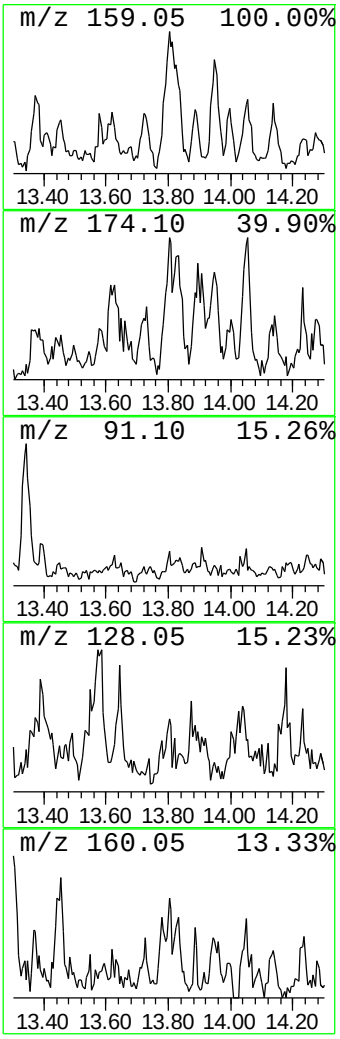
****

Peak Number 17 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	2.53 ng/ul	134497	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	87
2	1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	86
3	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	86
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	83
5	1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	78





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

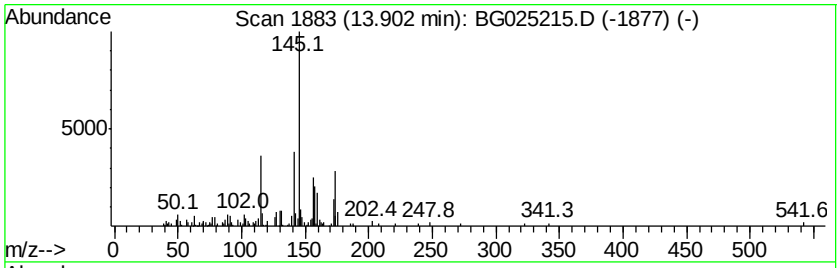
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

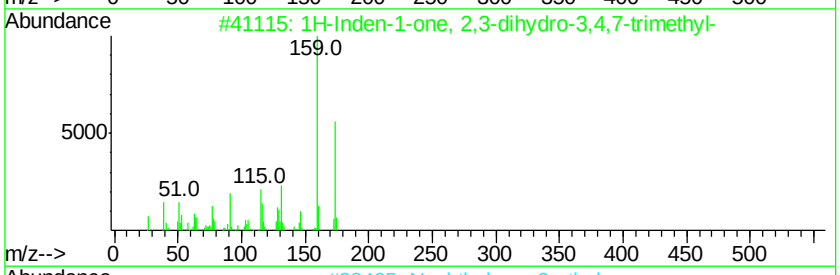
****

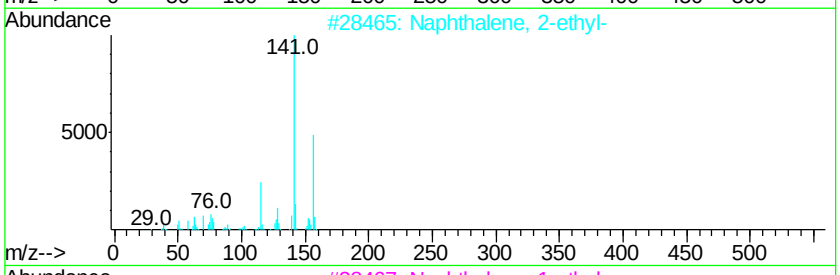
Peak Number 18 unknown-03 Concentration Rank 18

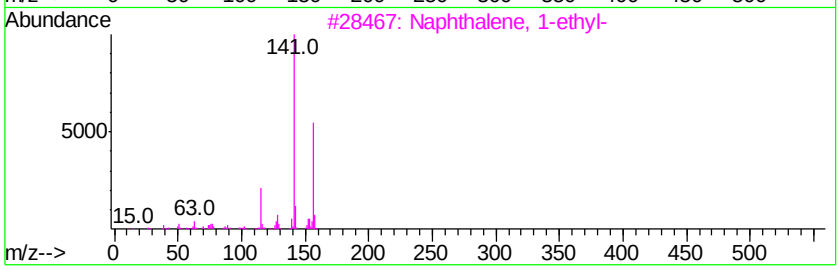
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	2.79 ng/ul	148325	Acenaphthene-d10	14.86

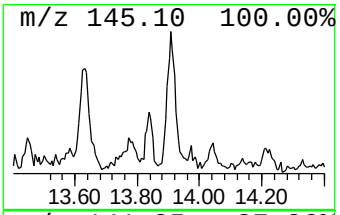
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	35
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	30
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	30
4	5-Amino-8-methoxyquinoline	174	C10H10N2O	070945-35-6	27
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	25

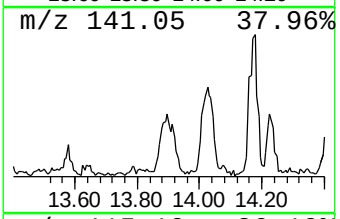


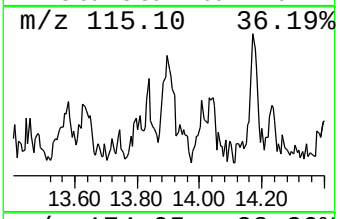


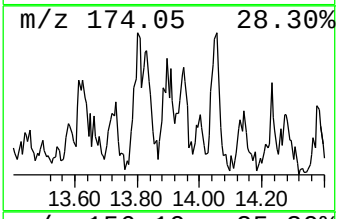


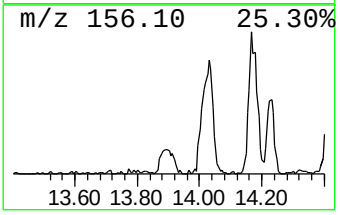












Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

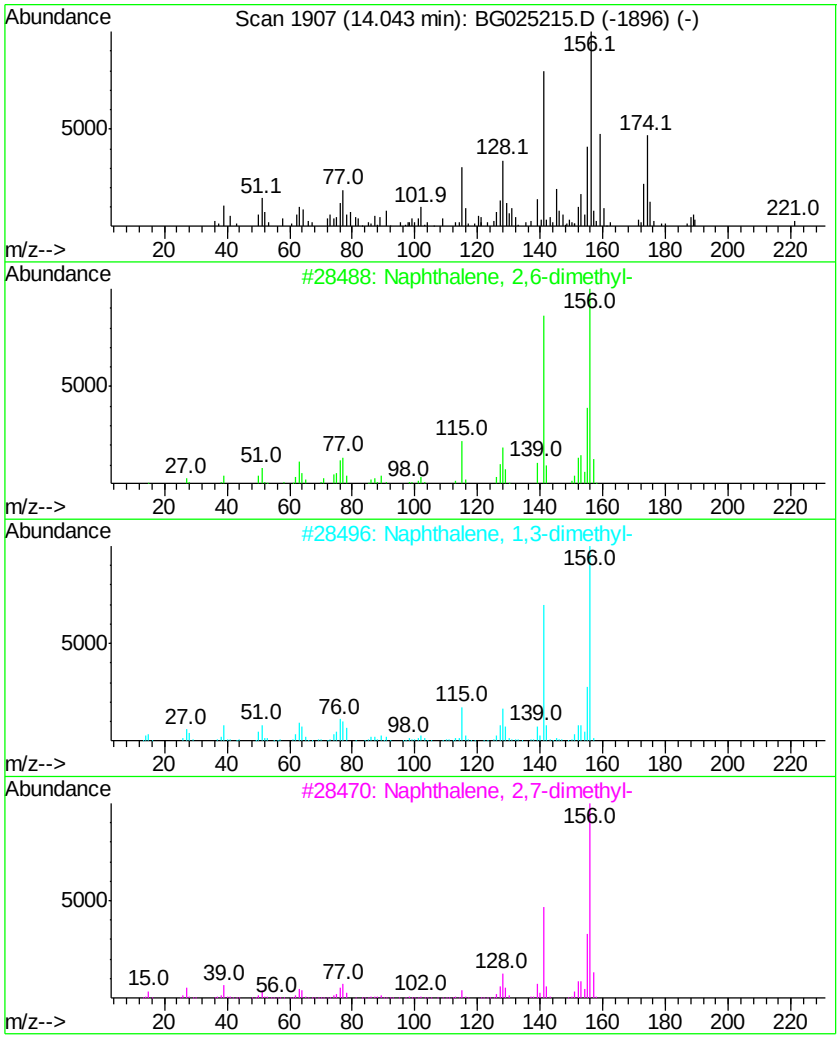
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

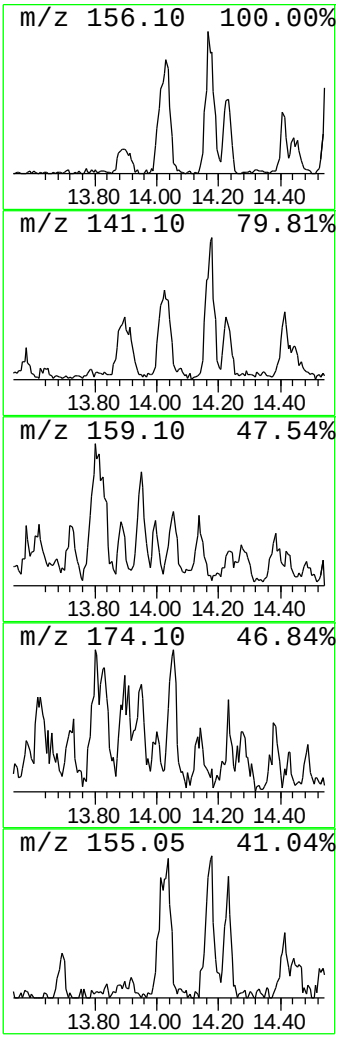
****

Peak Number 19 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	6.20 ng/ul	328984	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	89
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	89
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	76
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	68
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	64





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

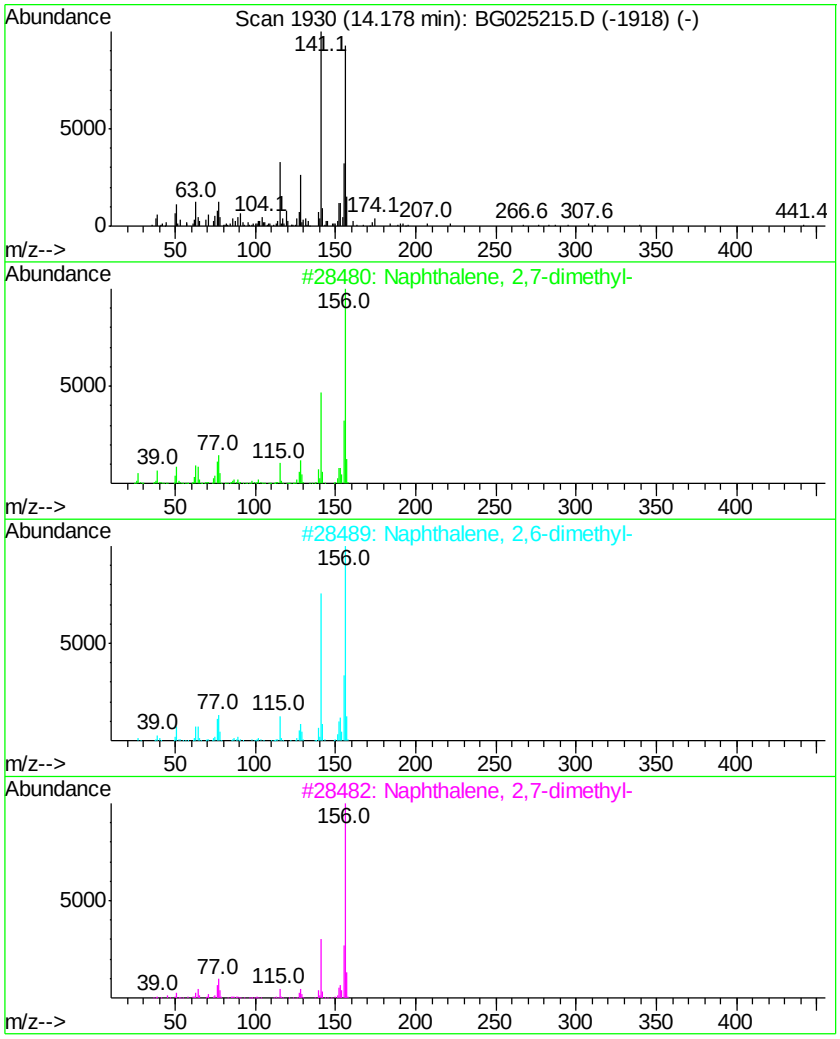
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

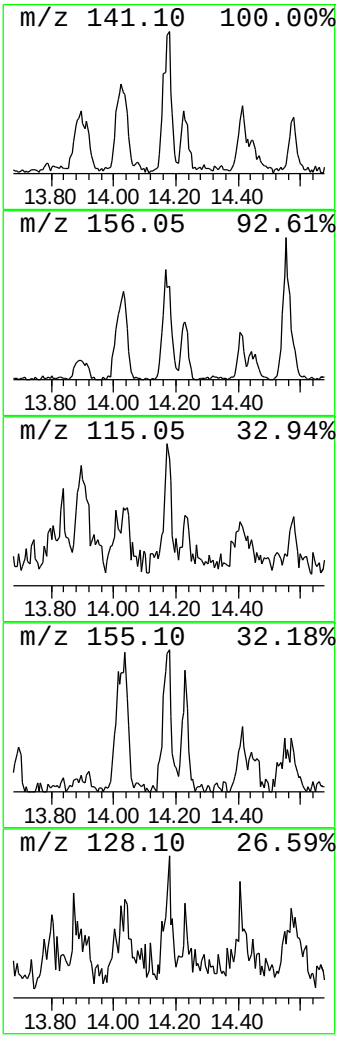
****

Peak Number 20 Naphthalene, 2,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	5.40 ng/ul	286942	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	91
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	87





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

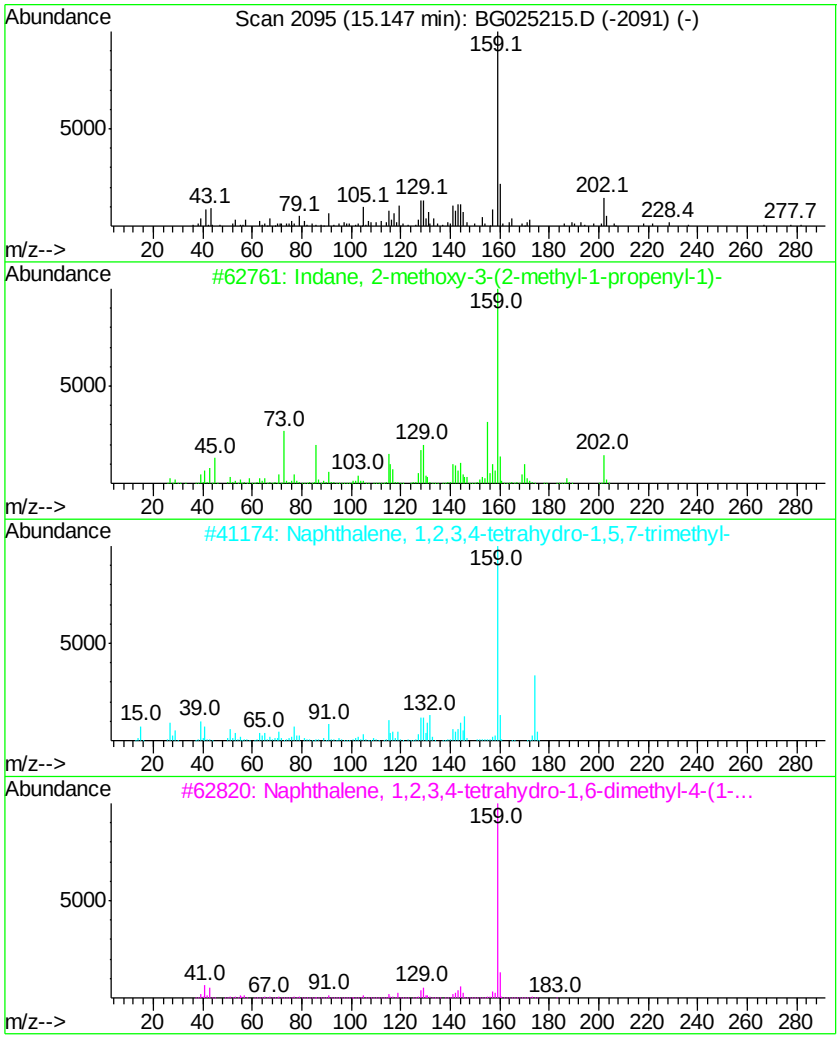
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

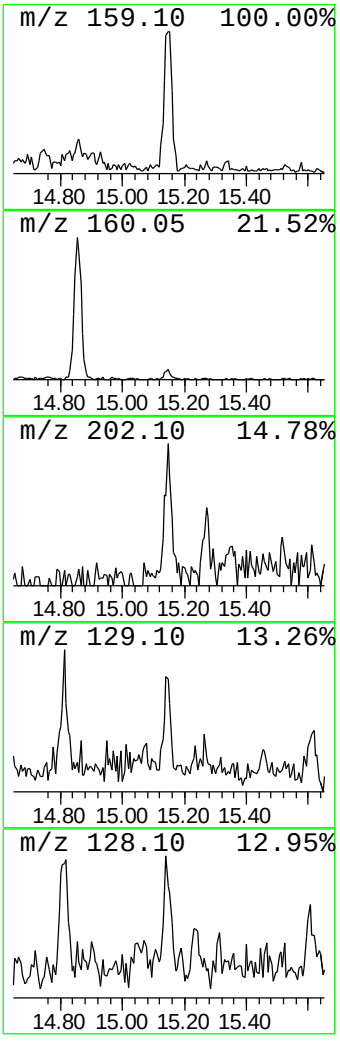
****

Peak Number 21 Indane, 2-methoxy-3-(2-meth... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	2.69 ng/ul	142808	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane, 2-methoxy-3-(2-methyl-1-...	202	C14H18O	1000196-88-7	74
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-55-0	72
3	Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	64
4	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	64
5	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	64





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

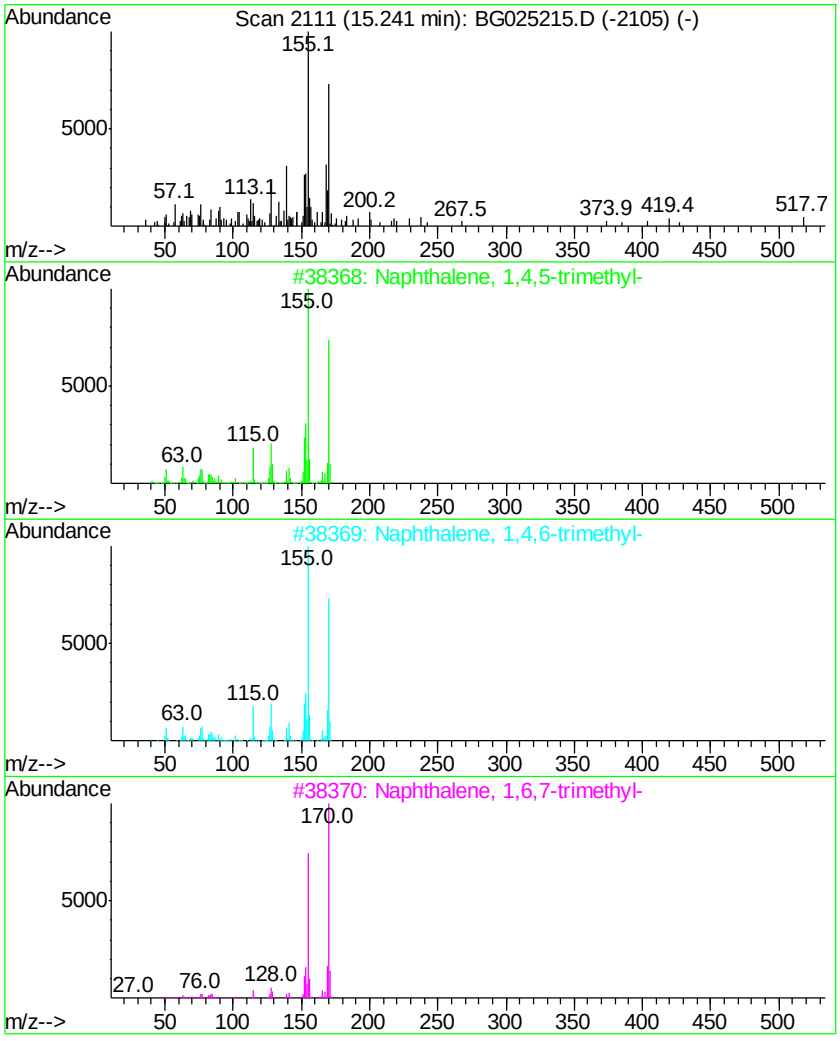
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

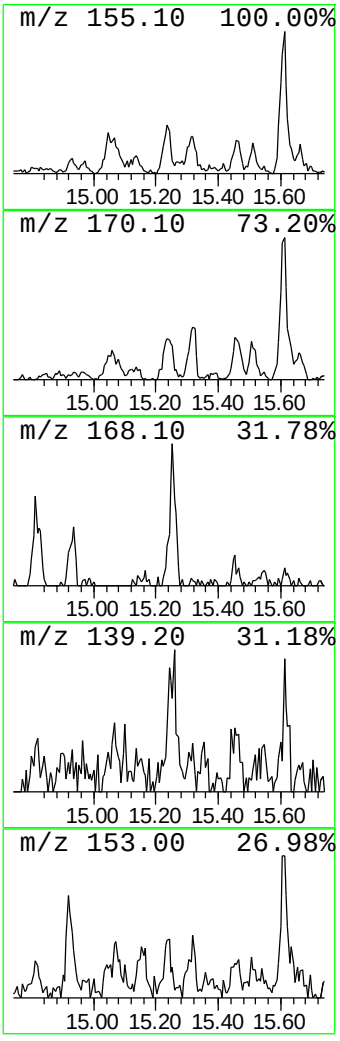
****

Peak Number 22 Naphthalene, 1,4,5-trimethyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	2.38 ng/ul	126563	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	90
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	90
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	81
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	81





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

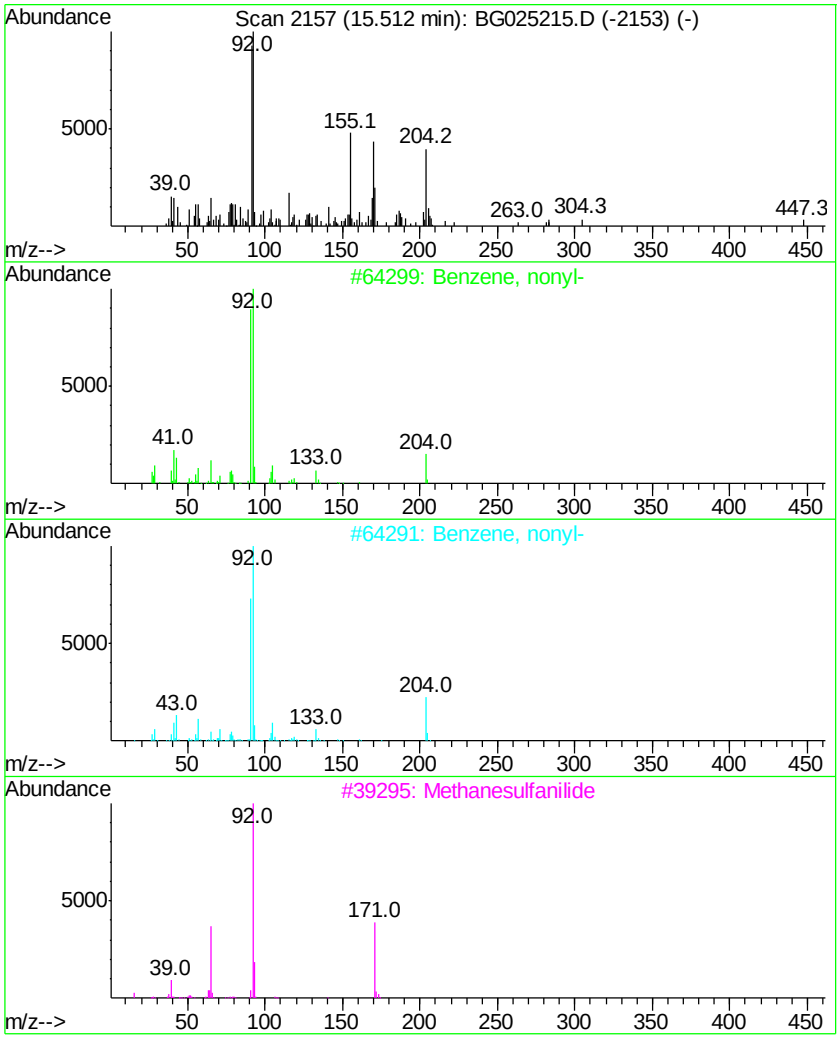
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

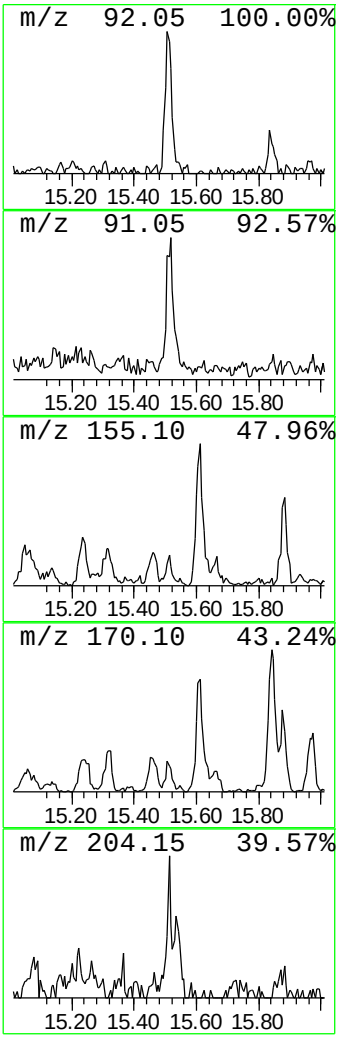
****

Peak Number 23 unknown-04 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	2.91 ng/ul	154492	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, nonyl-	204	C15H24	001081-77-2	45
2	Benzene, nonyl-	204	C15H24	001081-77-2	43
3	Methanesulfanilide	171	C7H9NO2S	001197-22-4	27
4	Benzene, (3-cyclopentylpropyl)-	188	C14H20	002883-12-7	25
5	Benzene, 1-methyl-4-(methylsulfo...	170	C8H10O2S	003185-99-7	25





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

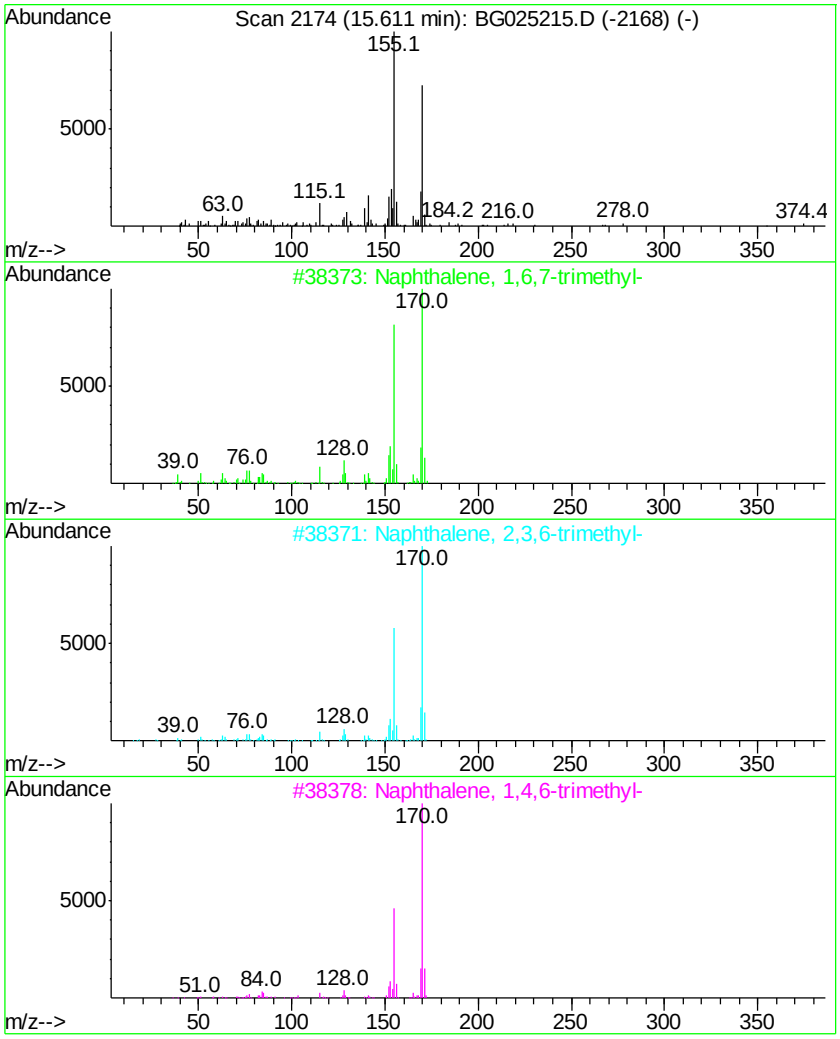
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

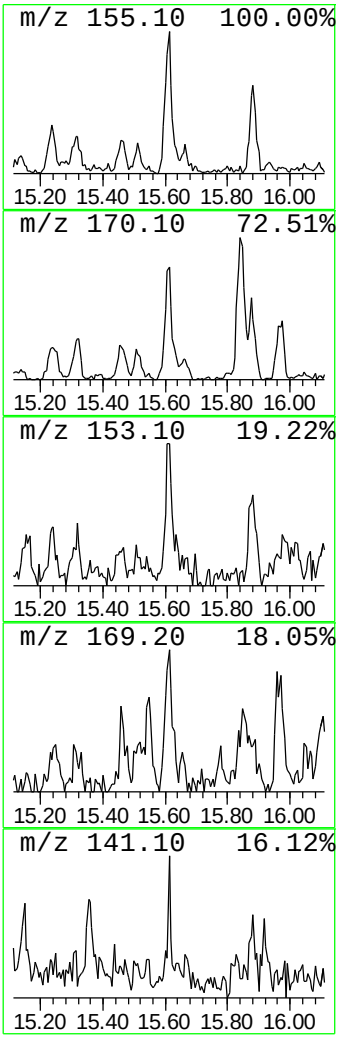
****

Peak Number 24 Naphthalene, 1,6,7-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	5.23 ng/ul	277499	Acenaphthene-d10	14.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	87
4	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	87
5	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	72





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

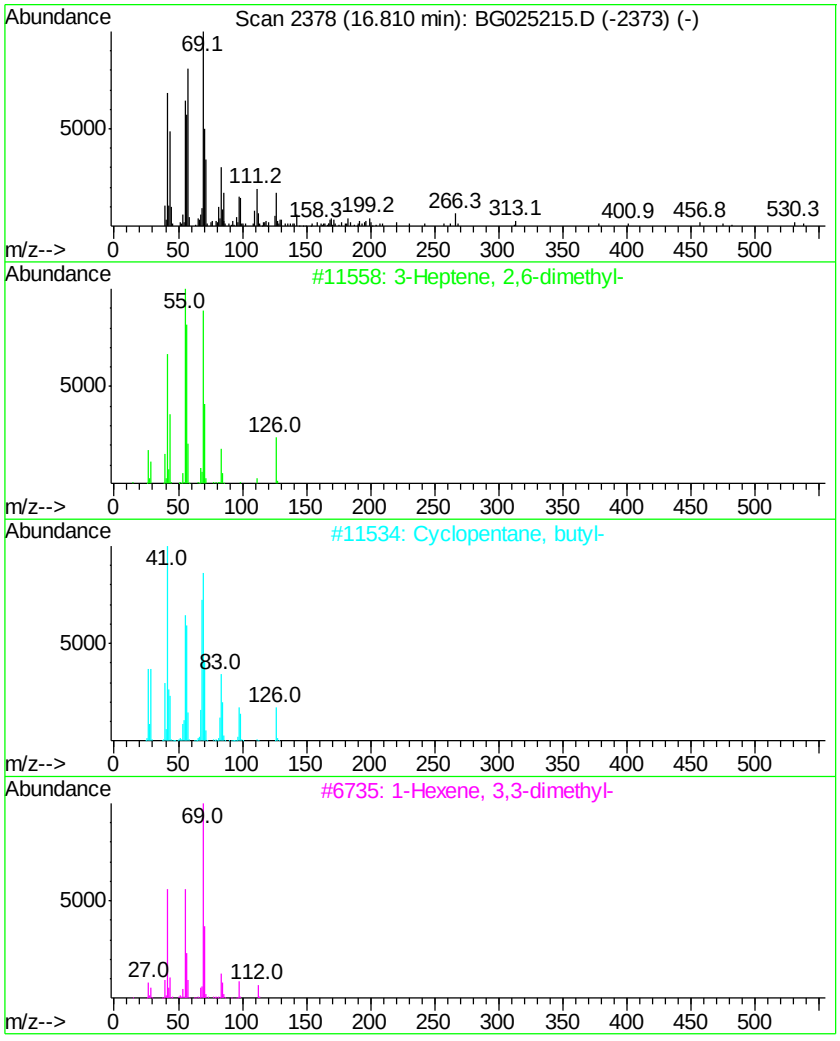
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

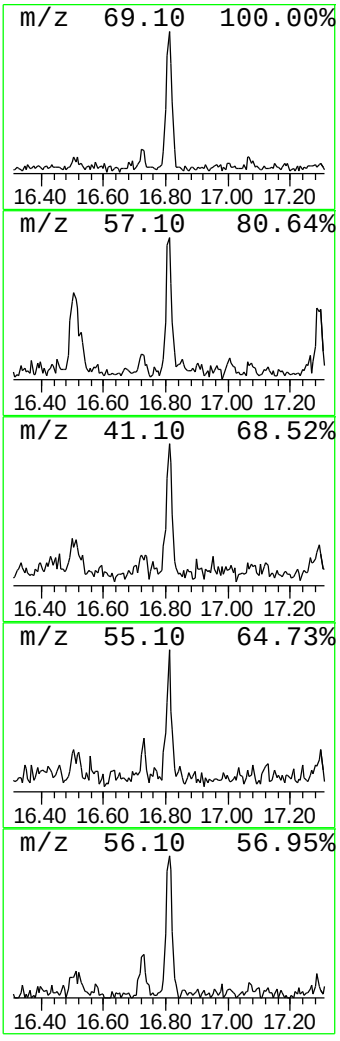
****

Peak Number 25 (DEL) Alkane: Cyclic16.81 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.81	2.53 ng/ul	170295	Phenanthrene-d10	17.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptene, 2,6-dimethyl-	126	C9H18	002738-18-3	72
2	Cyclopentane, butyl-	126	C9H18	002040-95-1	72
3	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	70
4	trans-2-Methyl-3-octene	126	C9H18	052937-36-7	58
5	Thieno[2,3-d]-1,3-thiaselenol-2-...	238	C5H2S3Se	081803-09-0	53





Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

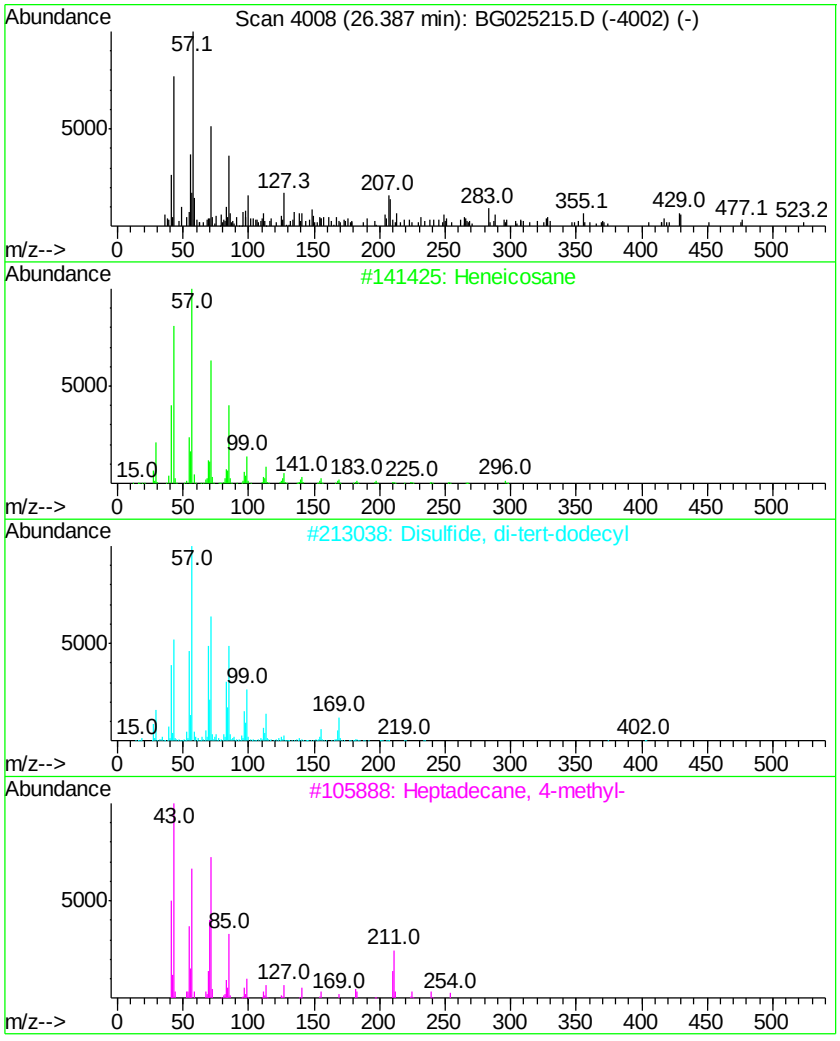
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

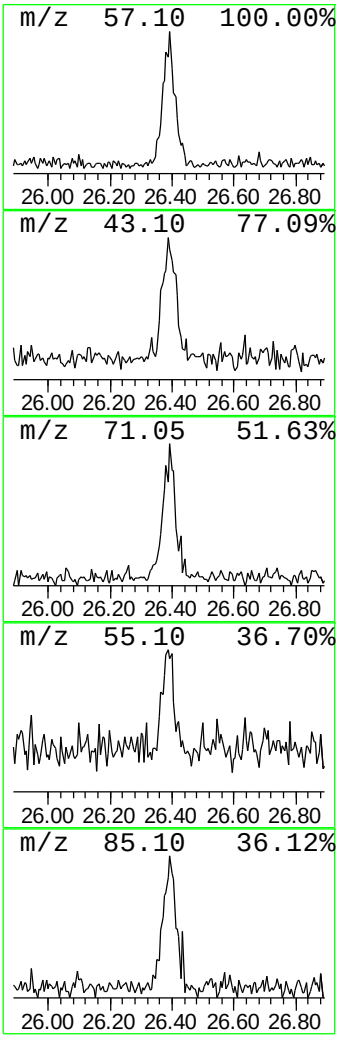
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.39	2.89 ng/ul	220400	Perylene-d12	25.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	64
2	Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	62
3	Heptadecane, 4-methyl-	254	C18H38	026429-11-8	58
4	Tricosane	324	C23H48	000638-67-5	58
5	Tritriacontane	465	C33H68	000630-05-7	58





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

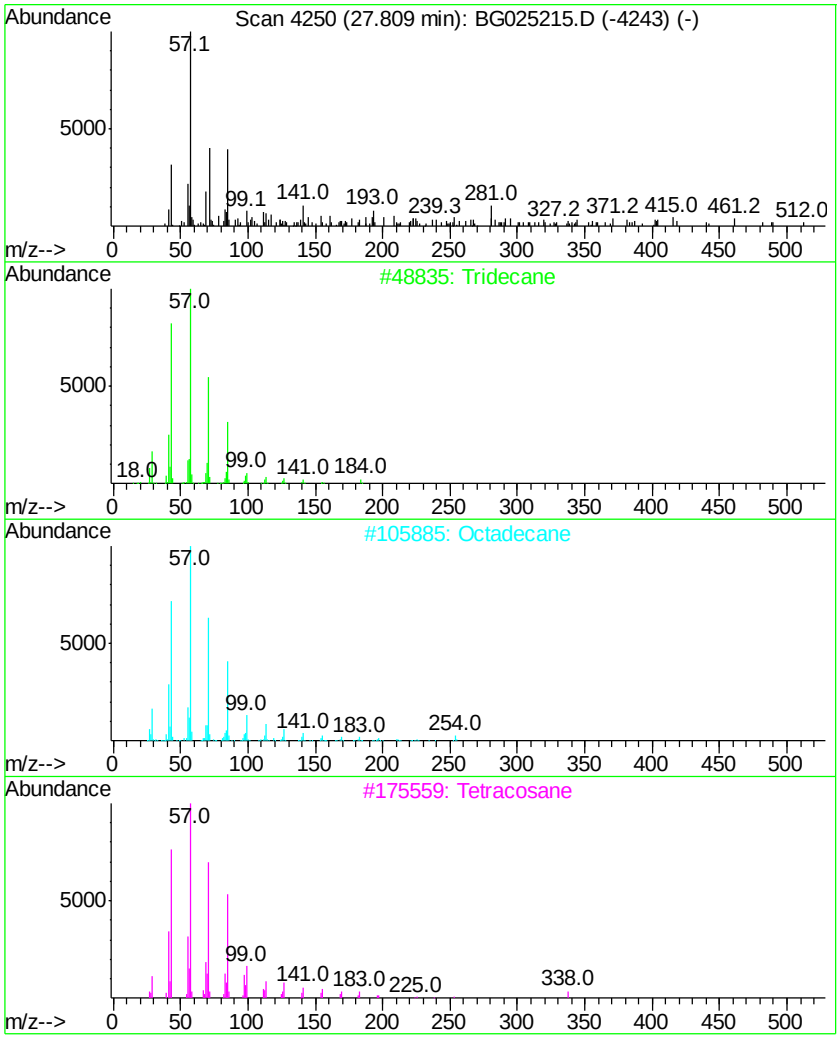
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

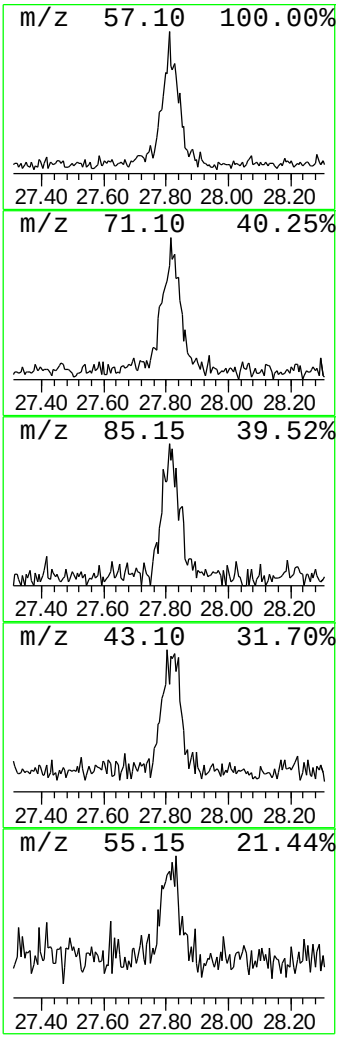
****

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.81	3.07 ng/ul	234078	Perylene-d12	25.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	58
2	Octadecane	254	C18H38	000593-45-3	58
3	Tetracosane	338	C24H50	000646-31-1	58
4	Tetracosane	338	C24H50	000646-31-1	58
5	Tridecane	184	C13H28	000629-50-5	52





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121516\
Data File : BG025215.D
Acq On : 15 Dec 2016 17:20
Operator : UM/SJ
Sample : H5995-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA869

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzofuran, 4,7-d...	11.29	7.0	ng/ul	219618	2	11.06	632137	20.0
1H-Indazole, 5,7-...	11.42	4.4	ng/ul	139521	2	11.06	632137	20.0
2-Propenal, 2-met...	11.46	9.8	ng/ul	311375	2	11.06	632137	20.0
4-Methyl-2H-benzo...	11.53	3.3	ng/ul	104688	2	11.06	632137	20.0
Benzene, hexyl-	12.00	2.7	ng/ul	85254	2	11.06	632137	20.0
unknown-01	12.09	3.2	ng/ul	102025	2	11.06	632137	20.0
1H-Indene, 4,7-di...	12.26	5.1	ng/ul	159920	2	11.06	632137	20.0
1H-Indene, 2,3-di...	12.36	2.9	ng/ul	91529	2	11.06	632137	20.0
Naphthalene, 1,2,...	12.57	2.5	ng/ul	79310	2	11.06	632137	20.0
1H-Benzimidazole,...	12.84	2.8	ng/ul	88887	2	11.06	632137	20.0
Benzocycloheptatr...	12.91	5.8	ng/ul	183917	2	11.06	632137	20.0
2,5,6-Trimethylbe...	13.00	3.6	ng/ul	190187	3	14.86	1061870	20.0
unknown-02	13.09	2.3	ng/ul	123853	3	14.86	1061870	20.0
1,2,3-Trimethylin...	13.58	2.8	ng/ul	147779	3	14.86	1061870	20.0
3-Ethyl-3-phenyl-...	13.63	2.1	ng/ul	111753	3	14.86	1061870	20.0
Naphthalene, 1,2,...	13.80	2.5	ng/ul	134497	3	14.86	1061870	20.0
unknown-03	13.90	2.8	ng/ul	148325	3	14.86	1061870	20.0
Naphthalene, 2,6-...	14.04	6.2	ng/ul	328984	3	14.86	1061870	20.0
Naphthalene, 2,7-...	14.18	5.4	ng/ul	286942	3	14.86	1061870	20.0
Indane, 2-methoxy...	15.15	2.7	ng/ul	142808	3	14.86	1061870	20.0
Naphthalene, 1,4,...	15.24	2.4	ng/ul	126563	3	14.86	1061870	20.0
unknown-04	15.51	2.9	ng/ul	154492	3	14.86	1061870	20.0
Naphthalene, 1,6,...	15.61	5.2	ng/ul	277499	3	14.86	1061870	20.0
(DEL) Alkane: Cyc...	16.81	2.5	ng/ul	170295	4	17.60	1344660	20.0
(DEL) Alkane: Str...	26.39	2.9	ng/ul	220400	6	25.31	1524950	20.0
(DEL) Alkane: Str...	27.81	3.1	ng/ul	234078	6	25.31	1524950	20.0