LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\
Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.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	3.416	69	73	78	rVB2	232471	329194	1.44%	0.135%
2	3.734	121	127	132	rVV	240848	369329	1.62%	0.152%
3	3.875	144	151	161	rVB5	137769	323046	1.41%	0.133%
4	4.181	194	203	211	rVV5	114768	304626	1.33%	0.125%
5	4.828	308	313	317	rBV	330988	470778	2.06%	0.193%
6	5.181	365	373	382	rBV	14963347	22846387	100.00%	9.386%
7	5.328	389	398	411	rVB	4814149	6835122	29.92%	2.808%
8	6.146	527	537	551	rBV	2177554	3376399	14.78%	1.387%
9	6.322	559	567	576	rVB2	207094	347924	1.52%	0.143%
10	6.628	613	619	630	rBV	4032853	6027965	26.38%	2.476%
11	6.740	632	638	642	rVV	858745	1234495	5.40%	0.507%
12	6.799	642	648	656	rVV2	2246165	3846462	16.84%	1.580%
13	6.875	656	661	672	rVB	1295657	2028847	8.88%	0.833%
14	6.981	672	679	684	rBV	1127631	1844261	8.07%	0.758%
15	7.040	684	689	701	rVB	3603945	5489985	24.03%	2.255%
16	7.175	707	712	719	rVV	1180388	1858492	8.13%	0.764%
17	7.293	726	732	740	rVB	7194673	11147781	48.79%	4.580%
18	7.510	764	769	772	rVV	252559	398995	1.75%	0.164%
19	7.546	772	775	781	rVV	411847	566157	2.48%	0.233%
20	7.646	785	792	795	rBV	1211779	1912342	8.37%	0.786%
21	7.681	795	798	806	rVV	645314	992281	4.34%	0.408%
22	7.775	806	814	821	rVB3	275813	609201	2.67%	0.250%
23	7.910	833	837	841	rVV3	176734	313959	1.37%	0.129%
24	8.016	848	855	868	rVB	6318947	9857651	43.15%	4.050%
25	8.134	868	875	879	rBV	1488122	2181244	9.55%	0.896%
26	8.187	879	884	893	rVB	1536884	2371823	10.38%	0.974%
27	8.281	893	900	905	rBV	1924368	3005822	13.16%	1.235%
28	8.351	905	912	921	rVV2	1021939	2206567	9.66%	0.907%
29	8.446	921	928	937	rVB	1000712	1683943	7.37%	0.692%
30	8.593	946	953	958	rBV	2094167	3120496	13.66%	1.282%
31	8.646	958	962	970	rVB	1367332	2064037	9.03%	0.848%
32	8.746	972	979	984	rVV	3957982	6085267	26.64%	2.500%
33	8.822	984	992	1002	rVV3	2268788	5639807	24.69%	2.317%
34	8.987	1012	1020	1027	rVB5	311358	733464	3.21%	0.301%

LSC Area Percent Report

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Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Title : SVOA CALIBRATION

35	9.075	1028	1035	1041	rBV3	188015	373342	1.63%	0.153%
36	9.263	1061	1067	1072	rVB	2041887	3031300	13.27%	1.245%
37	9.322	1072	1077	1085	rVB	2998960	4732420	20.71%	1.944%
38	9.516	1102	1110	1116	rBV2	1251848	2233826	9.78%	0.918%
39	9.628	1123	1129	1134	rBV4	376213	764509	3.35%	0.314%
40	9.681	1134	1138	1152	rVB	1242656	2134054	9.34%	0.877%
41	9.793	1152	1157	1158	rBV	201520	281945	1.23%	0.116%
42	9.834	1158	1164	1171	rVV3	4168488	7453211	32.62%	3.062%
43	9.904	1171	1176	1186	rVV2	390215	786533	3.44%	0.323%
44	10.016	1186	1195	1197	rVV5	311604	764202	3.34%	0.314%
45	10.057	1197	1202	1209	rVV2	1875500	3373762	14.77%	1.386%
46	10.134	1209	1215	1221	rVB	388524	653193	2.86%	0.268%
47	10.351	1248	1252	1255	rVV	180022	281582	1.23%	0.116%
48	10.428	1255	1265	1269	rVV2	1840063	3928742	17.20%	1.614%
49	10.481	1269	1274	1281	rVV	3182581	5705255	24.97%	2.344%
50	10.545	1281	1285	1289	rVV	470453	755112	3.31%	0.310%
51	10.598	1289	1294	1298	rVV	381723	611949	2.68%	0.251%
52	10.645	1298	1302	1307	rVB	272397	428359	1.87%	0.176%
53	10.710	1307	1313	1321	rVB2	217611	511699	2.24%	0.210%
54	11.087	1373	1377	1383	rVB	338270	529172	2.32%	0.217%
55	11.345	1415	1421	1427	rVB	475542	800433	3.50%	0.329%
56	11.540	1448	1454	1459	rBV	459852	730088	3.20%	0.300%
57	11.622	1464	1468	1477	rVB3	159043	291522	1.28%	0.120%
58	11.822	1496	1502	1511	rVB2	334703	623444	2.73%	0.256%
59	11.987	1519	1530	1539	rVB6	391779	938321	4.11%	0.385%
60	12.098	1544	1549	1557	rVB	736825	1191621	5.22%	0.490%
61	12.322	1574	1587	1592	rBV	2554895	4348478	19.03%	1.786%
62	12.398	1592	1600	1604	rVV	446558	1365270	5.98%	0.561%
63	12.457	1604	1610	1616	rVB	737048	1475857	6.46%	0.606%
64	13.069	1704	1714	1719	rVV	534042	1049825	4.60%	0.431%
65	13.122	1719	1723	1727	rVV	353163	536503	2.35%	0.220%
66	13.163	1727	1730	1736	rVB3	217225	331885	1.45%	0.136%
67	13.222	1736	1740	1752	rBV4	207872	464146	2.03%	0.191%
68	13.322	1752	1757	1760	rBV	202773	334691	1.46%	0.137%
69	13.469	1771	1782	1789	rBV6	515421	1749883	7.66%	0.719%
70	13.528	1789	1792	1800	rVV3	502717	923860	4.04%	0.380%
71	13.616	1800	1807	1812	rVV	951770	1836387	8.04%	0.754%

LSC Area Percent Report

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Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Title : SVOA CALIBRATION

72	13.669	1812	1816	1819	rVV2	730805	1173612	5.14%	0.482%
73	13.710	1819	1823	1833	rVV	3015389	4350175	19.04%	1.787%
74	13.857	1841	1848	1851	rVV	346325	652801	2.86%	0.268%
75	13.934	1857	1861	1866	rBV	504798	772676	3.38%	0.317%
76	13.992	1866	1871	1880	rVV	3330123	5514070	24.14%	2.265%
77	14.298	1915	1923	1930	rBV2	2447618	4102824	17.96%	1.686%
78	14.504	1948	1958	1969	rVV3	1616352	3671085	16.07%	1.508%
79	14.704	1987	1992	2001	rVV2	326359	546774	2.39%	0.225%
80	14.816	2001	2011	2024	rVV	1276931	2812562	12.31%	1.155%
81	14.928	2025	2030	2036	rVV5	232553	610956	2.67%	0.251%
82	15.192	2068	2075	2078	rVV2	142219	317379	1.39%	0.130%
83	15.239	2079	2083	2088	rVV	634812	943926	4.13%	0.388%
84	15.298	2088	2093	2098	rVV	4058002	5772683	25.27%	2.372%
85	15.428	2110	2115	2120	rBV	1164383	1573446	6.89%	0.646%
86	15.663	2150	2155	2161	rBV4	200430	341782	1.50%	0.140%
87	15.851	2183	2187	2202	rVB	352128	727913	3.19%	0.299%
88	16.051	2214	2221	2230	rBV4	199582	402067	1.76%	0.165%
89	17.051	2386	2391	2397	rBV2	2872885	4274626	18.71%	1.756%
90	17.151	2403	2408	2421	rVB2	4366917	6214033	27.20%	2.553%
91	17.963	2542	2546	2557	rVB5	128885	287306	1.26%	0.118%
92	19.463	2795	2801	2813	rBV	5380197	7217485	31.59%	2.965%
93	21.274	3103	3109	3123	rBV	3409474	4687555	20.52%	1.926%
94	22.333	3286	3289	3295	rVB	308246	367335	1.61%	0.151%
95	22.839	3371	3375	3382	rVB	414660	553907	2.42%	0.228%
96	23.374	3458	3466	3482	rBV2	3763773	7869310	34.44%	3.233%
97	23.510	3483	3489	3504	rVB2	2426315	4743398	20.76%	1.949%
98	24.045	3575	3580	3587	rVB	443797	788033	3.45%	0.324%
99	24.786	3700	3706	3713	rVB	362859	780790	3.42%	0.321%
100	25.645	3847	3852	3863	rVB2	242426	592330	2.59%	0.243%

Sum of corrected areas: 243413369

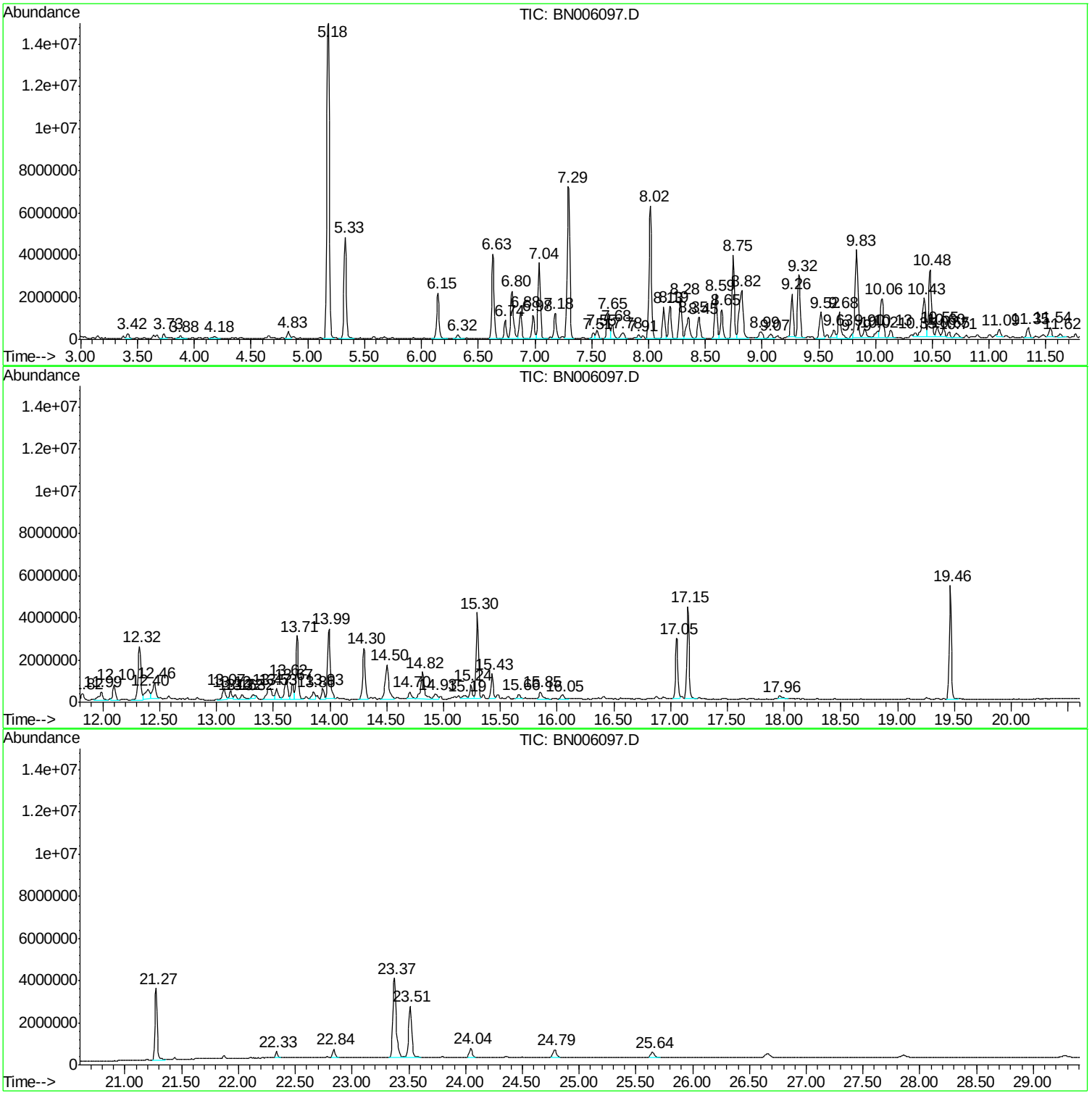
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\
Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\
Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

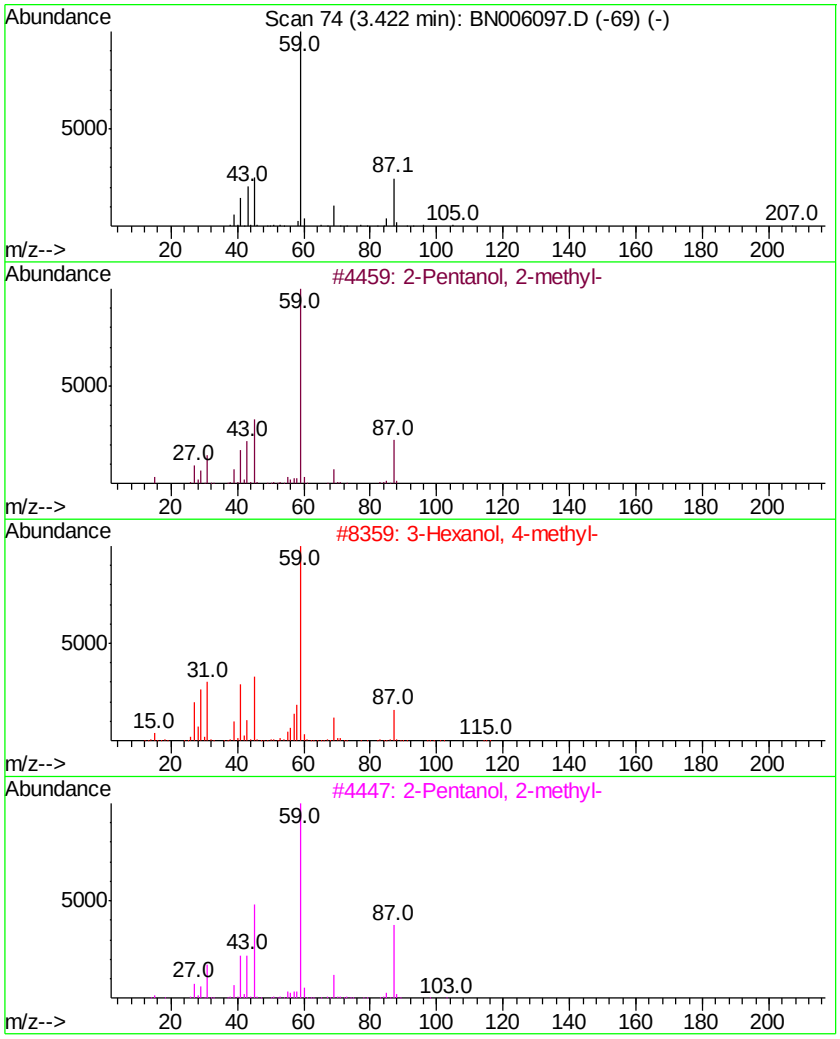
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

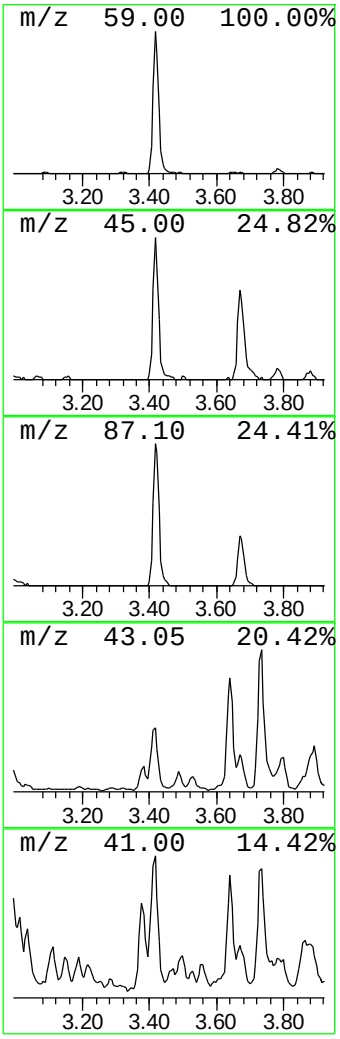
****

Peak Number 1 2-Pentanol, 2-methyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.42	3.44 ng/ul	329194	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	83
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	64
3	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	64
4	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	56
5	2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\
Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

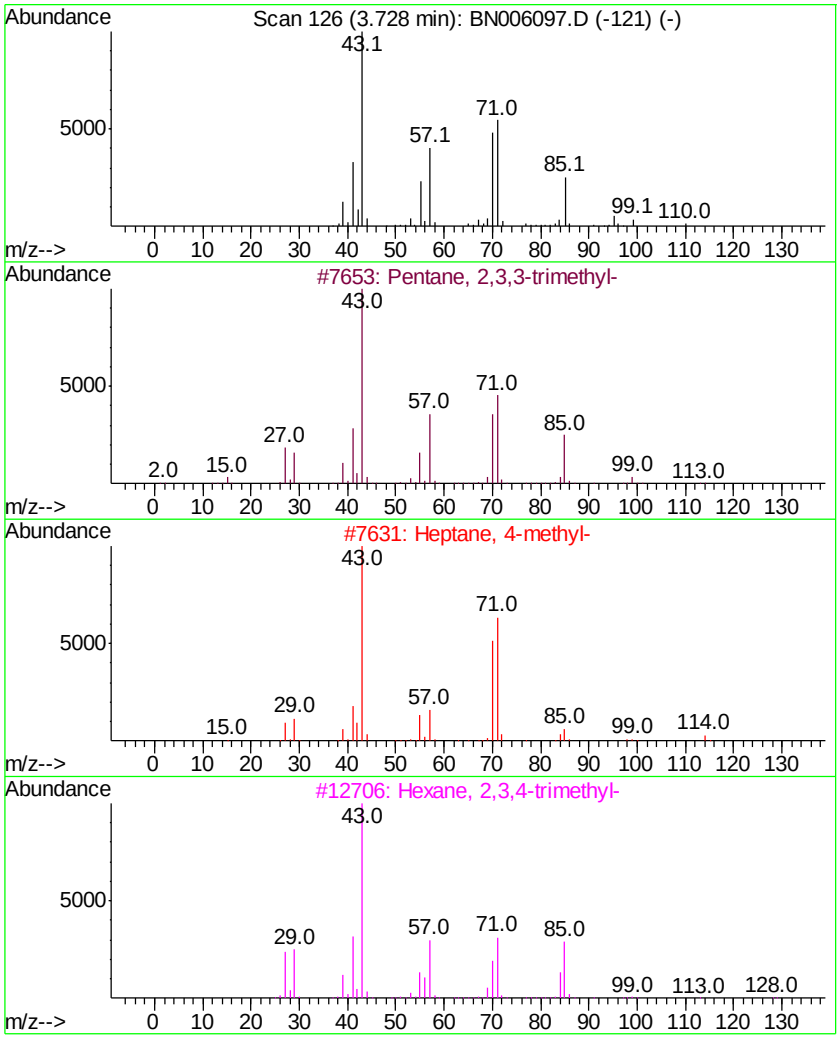
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

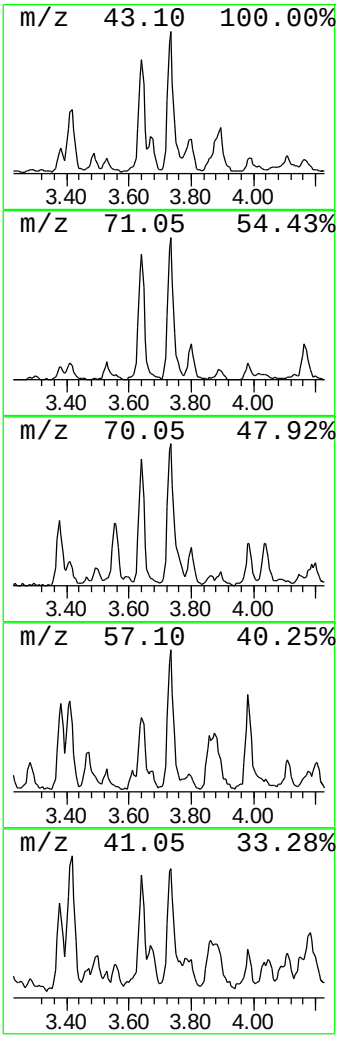
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	3.86 ng/ul	369329	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	83
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	59
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59





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Instrument : BNA_N
ClientSampleId : C0C52

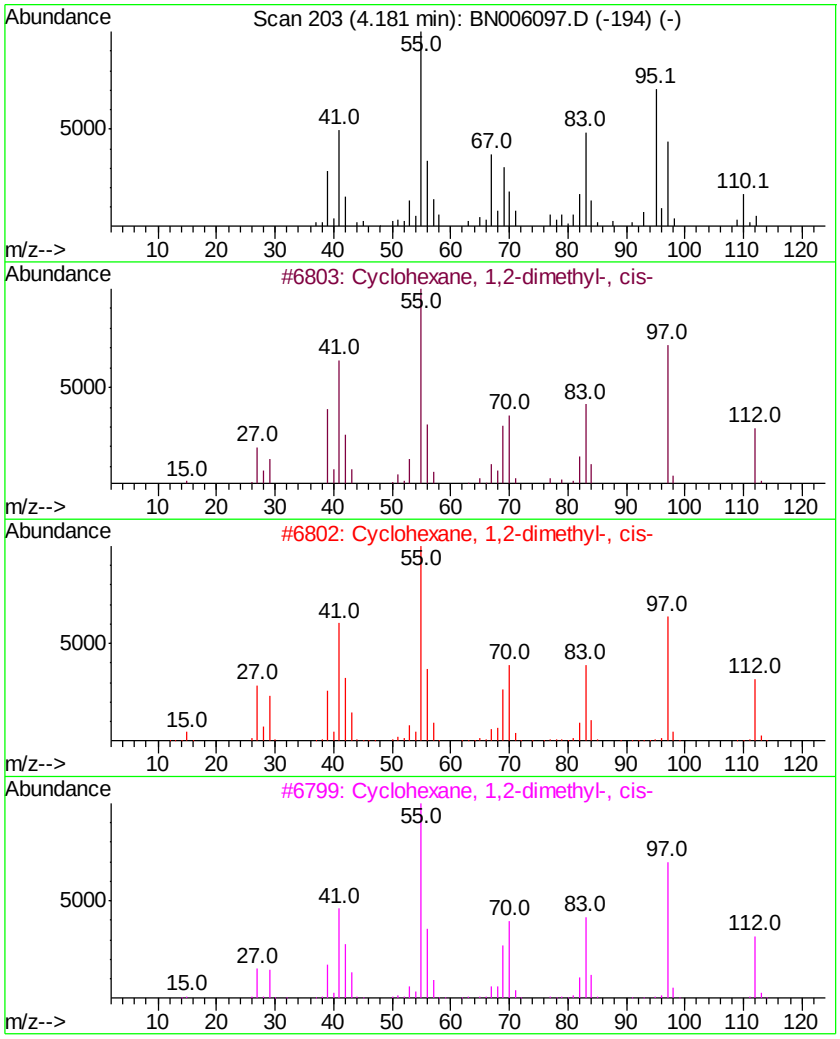
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

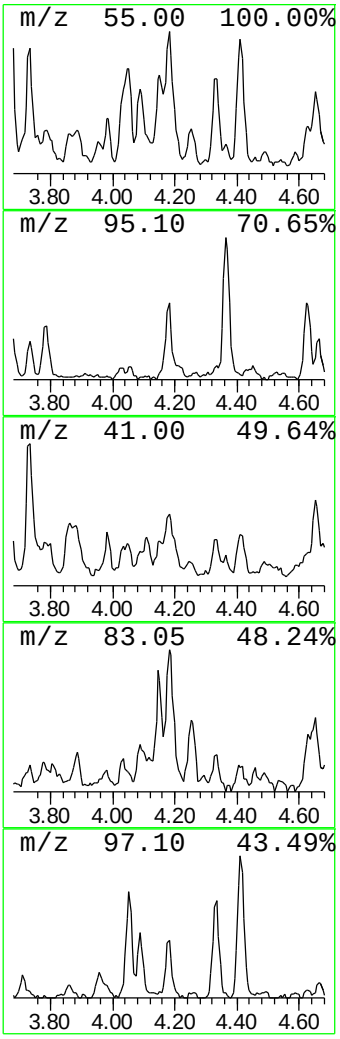
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Cyclic4.18 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.18	3.19 ng/ul	304626	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	58
2	Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	53
3	Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	49
4	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	43
5	1,4-Hexadiene, 2,3-dimethyl-	110	C8H14	018669-52-8	38





Library Search Compound Report

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Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

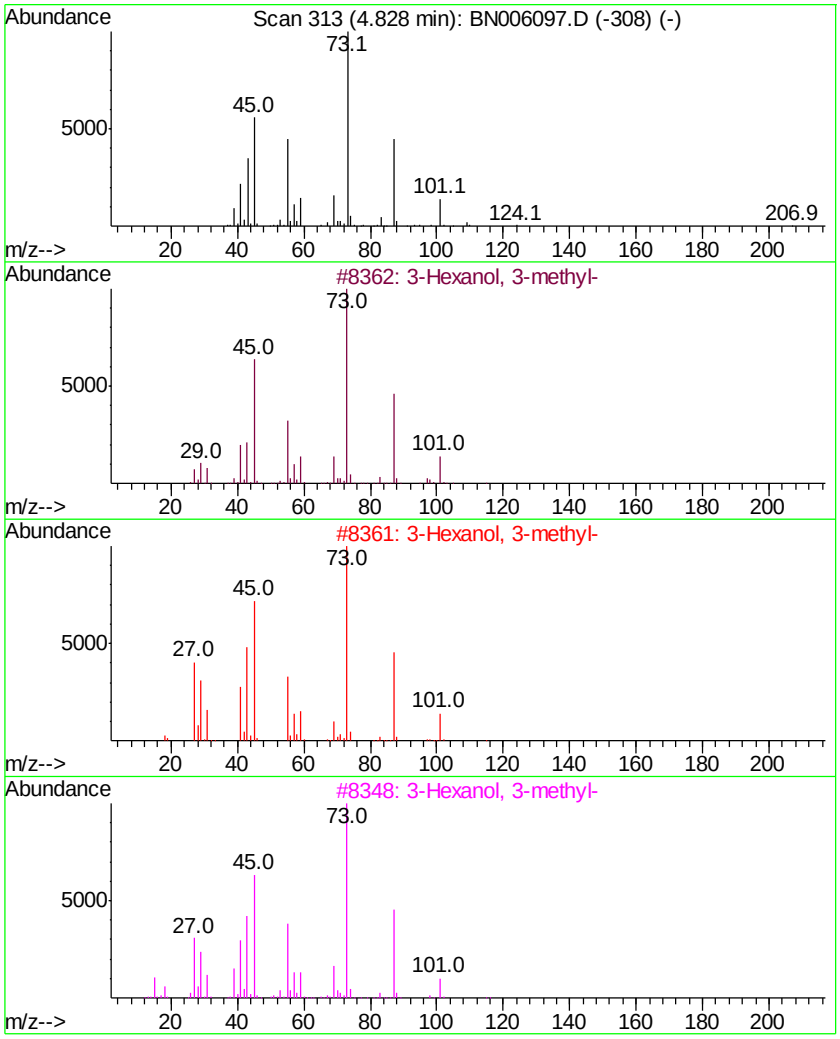
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

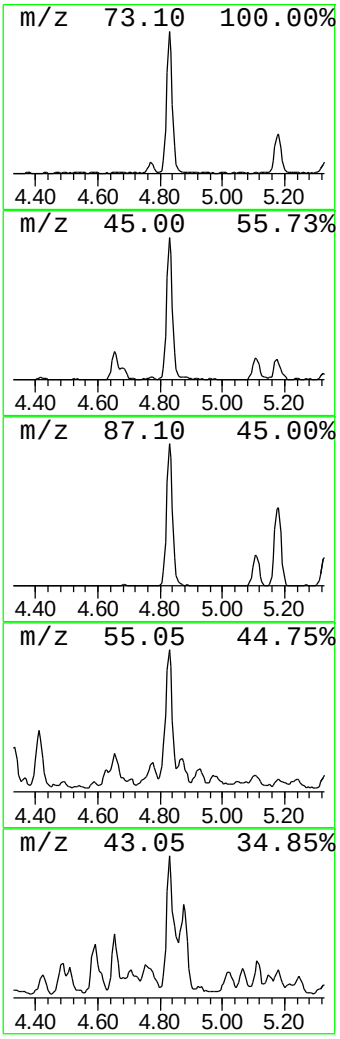
****

Peak Number 5 3-Hexanol, 3-methyl- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	4.92 ng/ul	470778	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	90
2	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	90
3	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	86
4	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	53
5	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	50





Library Search Compound Report

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Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

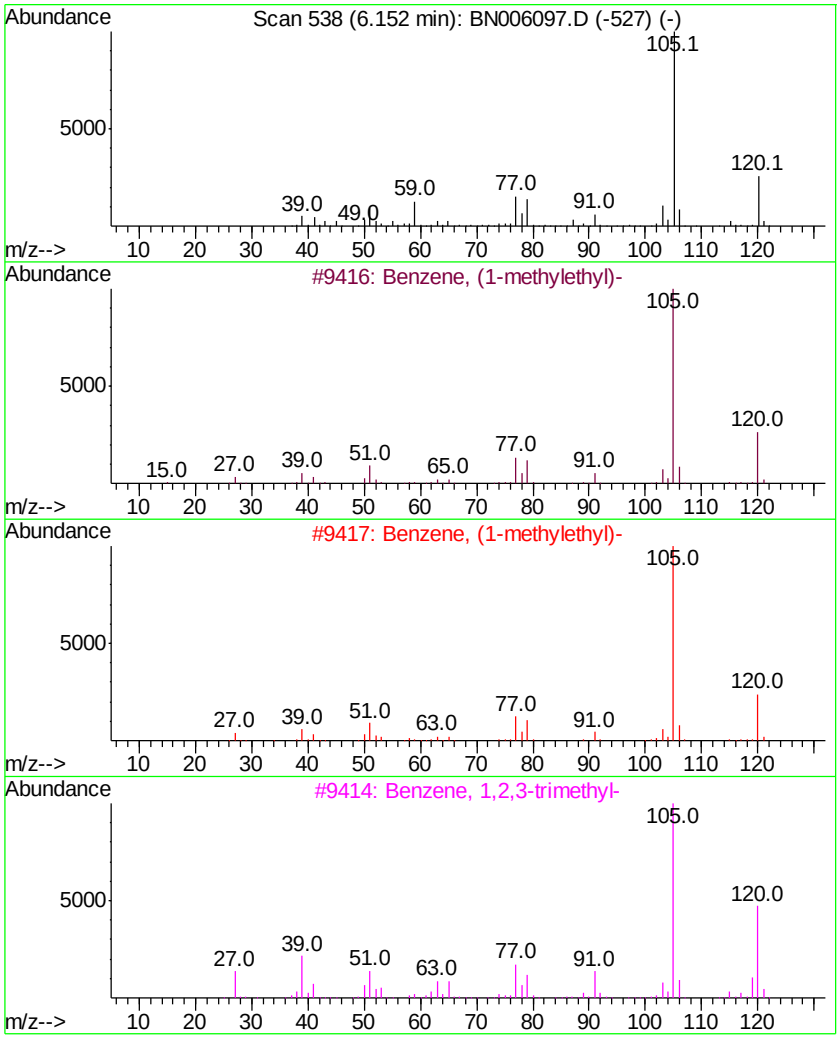
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

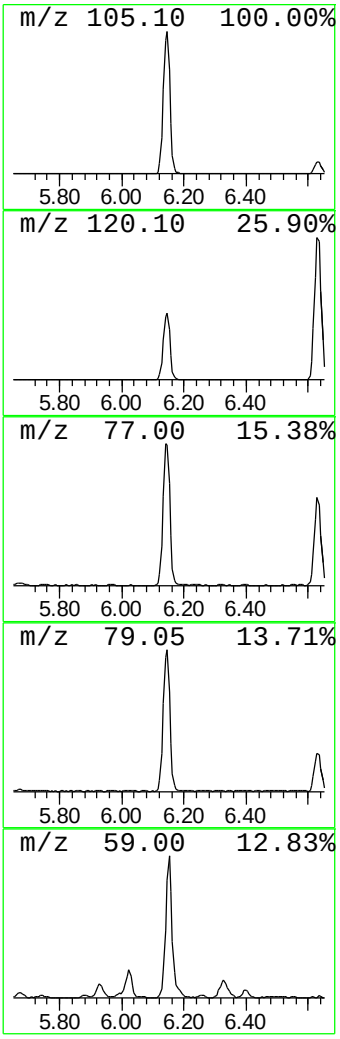
****

Peak Number 8 Benzene, (1-methylethyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.15	35.31 ng/ul	3376400	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	81





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Sample : K3045-08
Misc :
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Instrument : BNA_N
ClientSampleId : C0C52

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Quant Title : SVOA CALIBRATION

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

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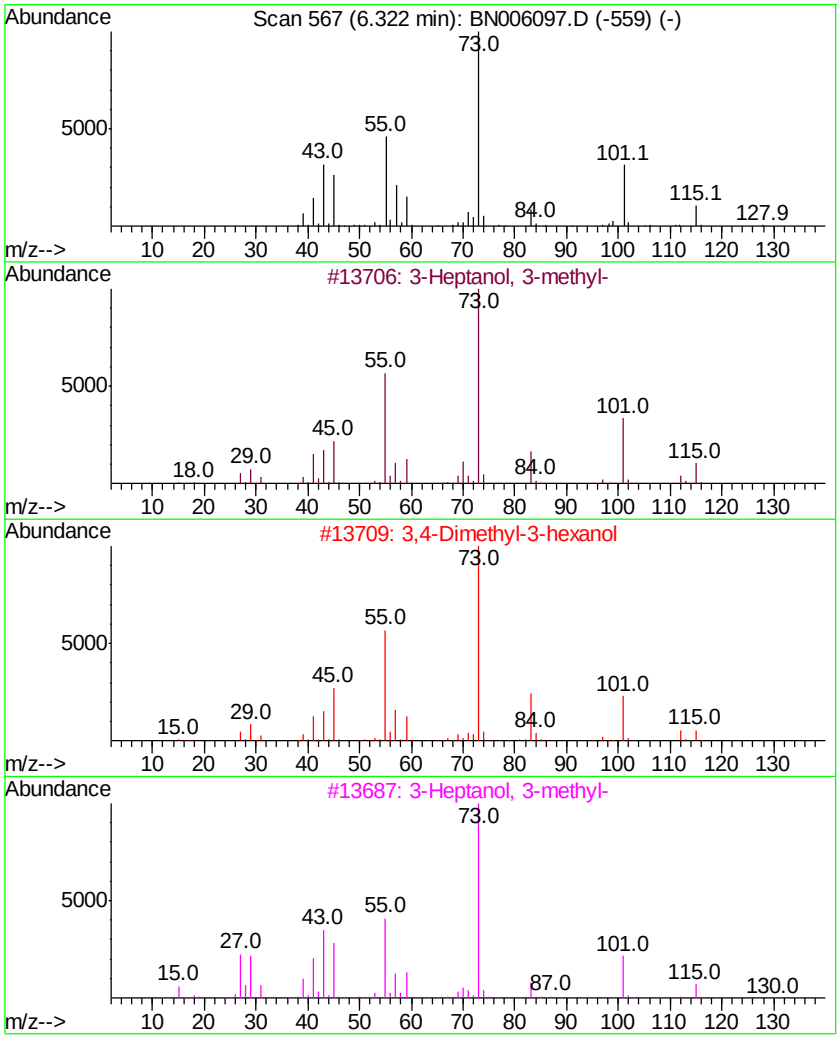
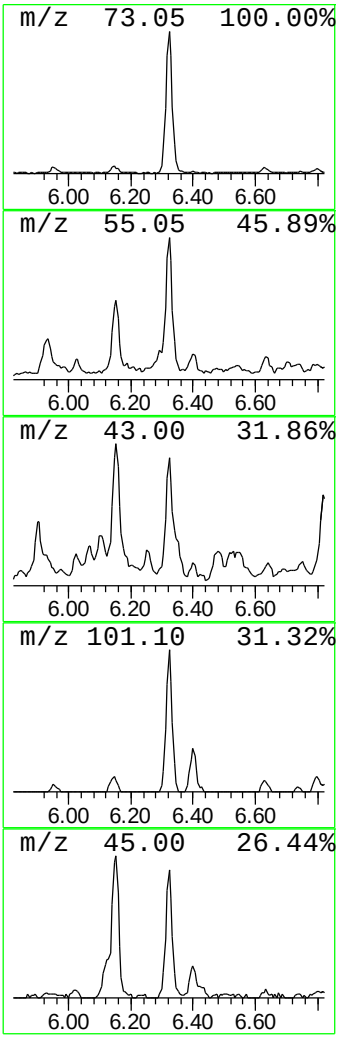
Peak Number 9 3,4-Dimethyl-3-hexanol Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	3.64 ng/ul	347924	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanol, 3-methyl-	130	C8H18O	005582-82-1	78
2	3,4-Dimethyl-3-hexanol	130	C8H18O	019550-08-4	72
3	3-Heptanol, 3-methyl-	130	C8H18O	005582-82-1	64
4	3,4-Dimethyl-3-hexanol	130	C8H18O	019550-08-4	56
5	3-Hexanol, 3,5-dimethyl-	130	C8H18O	004209-91-0	42





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ClientSampleId : C0C52

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Quant Title : SVOA CALIBRATION

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

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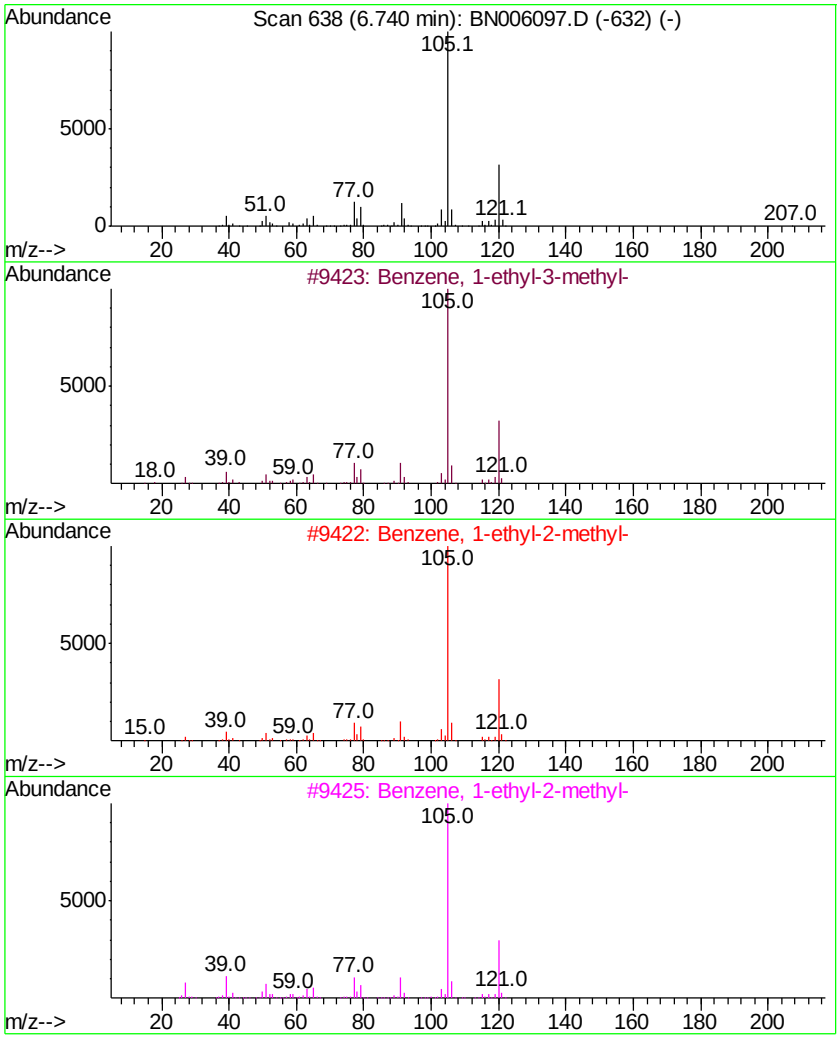
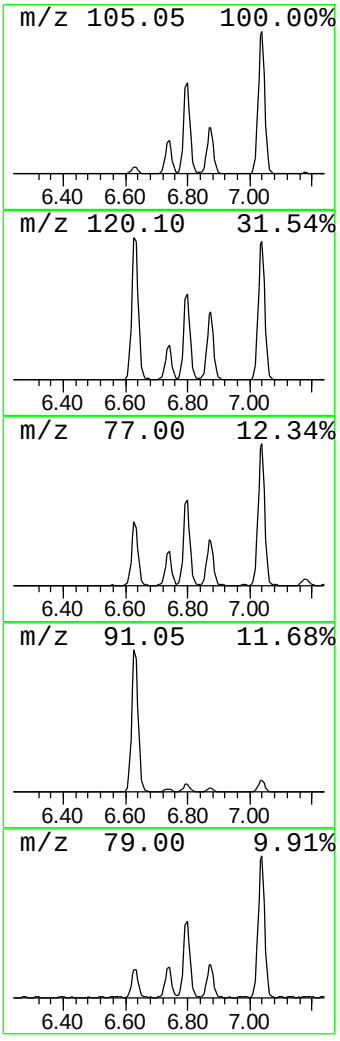
Peak Number 11 Benzene, 1-ethyl-3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	12.91 ng/ul	1234500	1,4-Dichlorobenzene-d4	7.65

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-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Instrument : BNA_N
ClientSampleId : C0C52

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Quant Title : SVOA CALIBRATION

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

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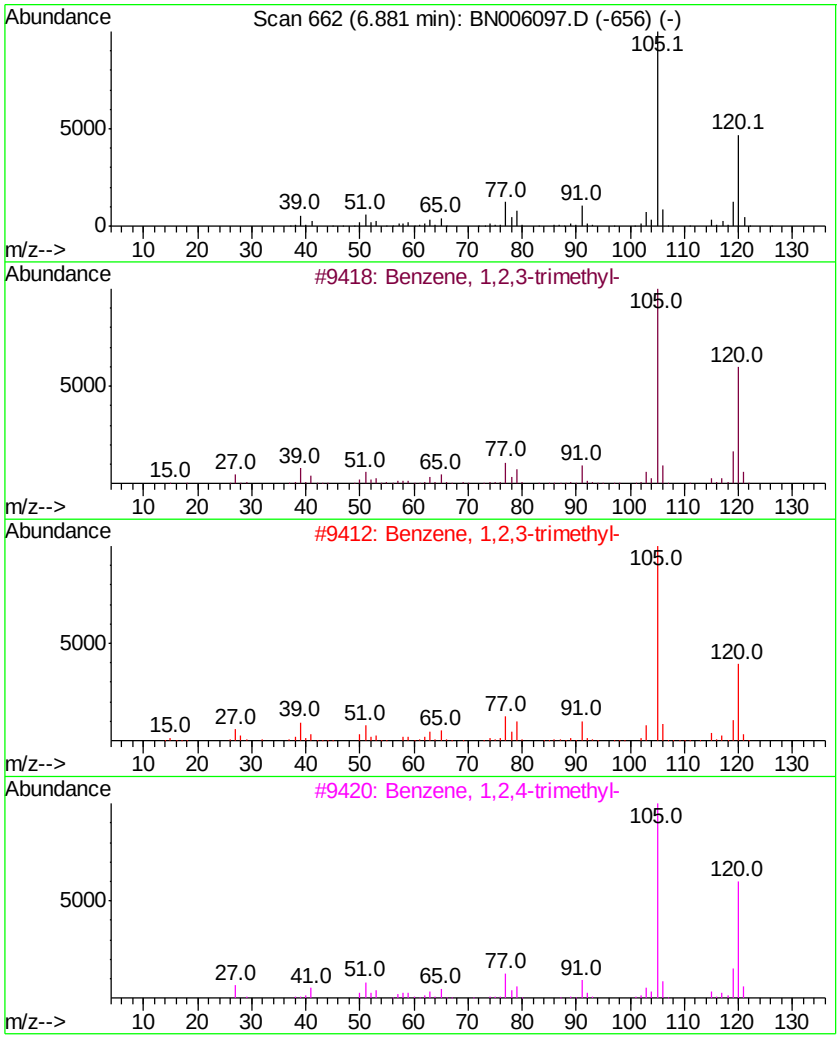
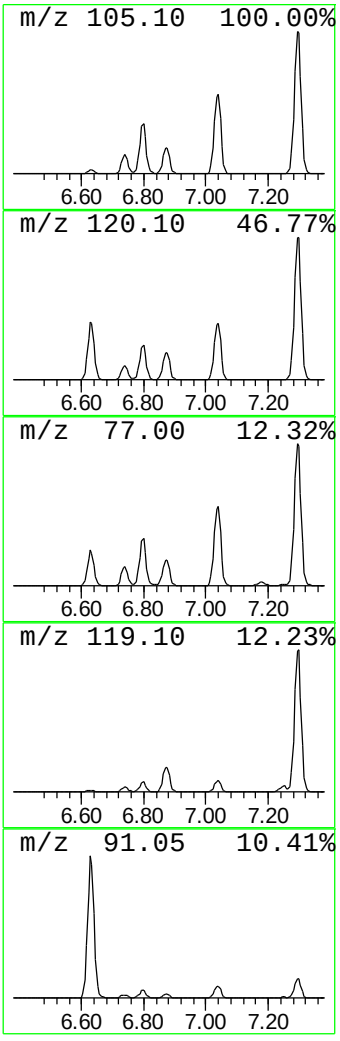
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	21.22 ng/ul	2028850	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Mesitylene	120	C9H12	000108-67-8	93
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93





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Instrument : BNA_N
ClientSampleId : C0C52

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Quant Title : SVOA CALIBRATION

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

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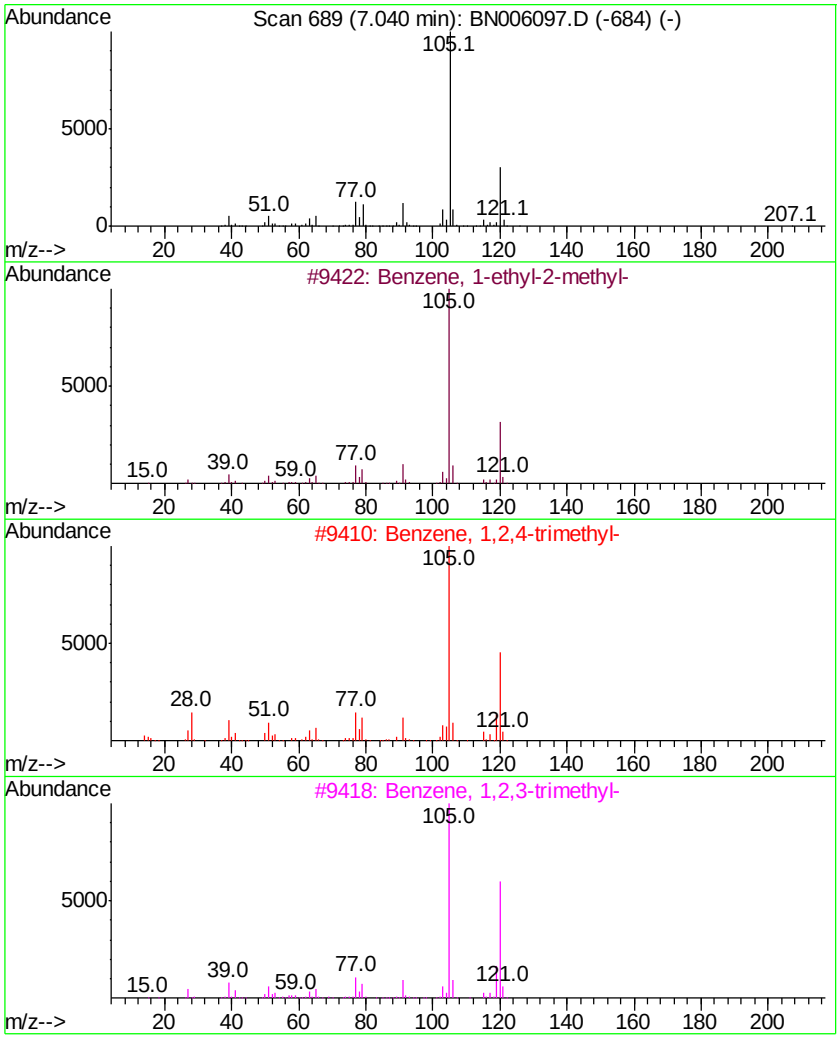
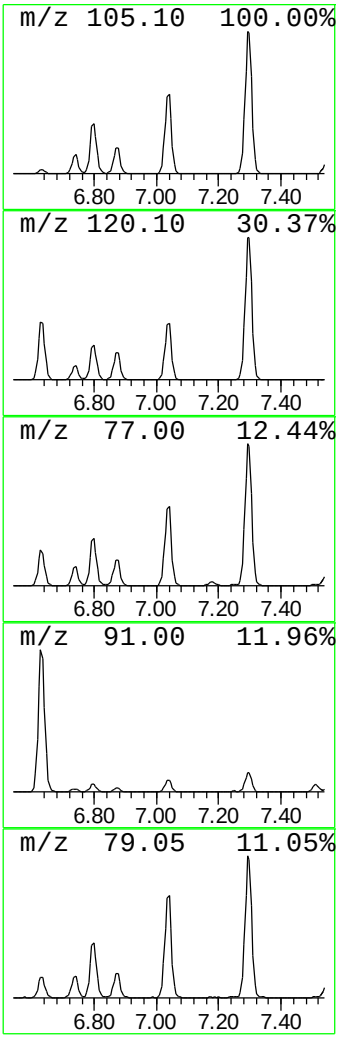
Peak Number 13 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	57.42 ng/ul	5489990	1,4-Dichlorobenzene-d4	7.65

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





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Instrument : BNA_N
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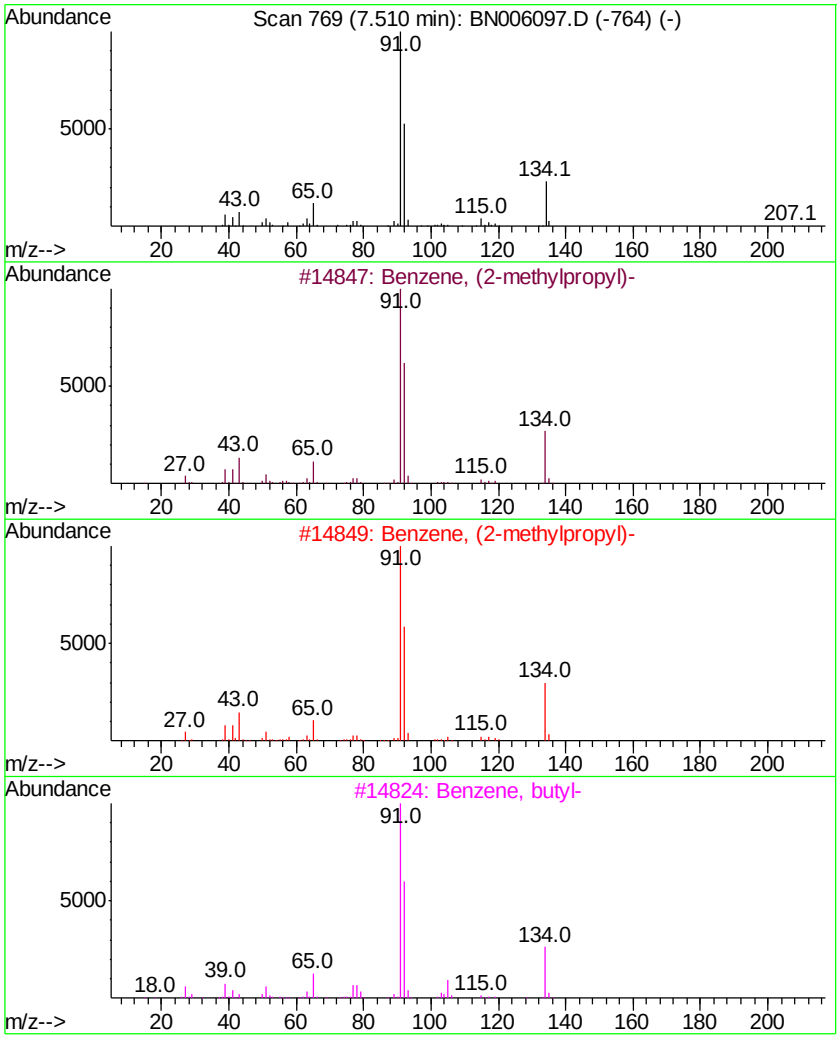
Quant Title : SVOA CALIBRATION

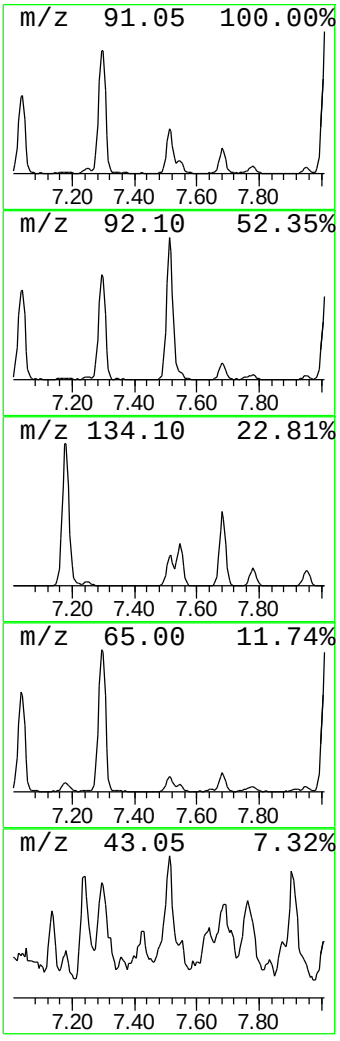
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, (2-methylpropyl)- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.51	4.17 ng/ul	398995	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
2	Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	90
3	Benzene, butyl-	134	C10H14	000104-51-8	80
4	Benzene, butyl-	134	C10H14	000104-51-8	80
5	Benzene, butyl-	134	C10H14	000104-51-8	72





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Instrument : BNA_N
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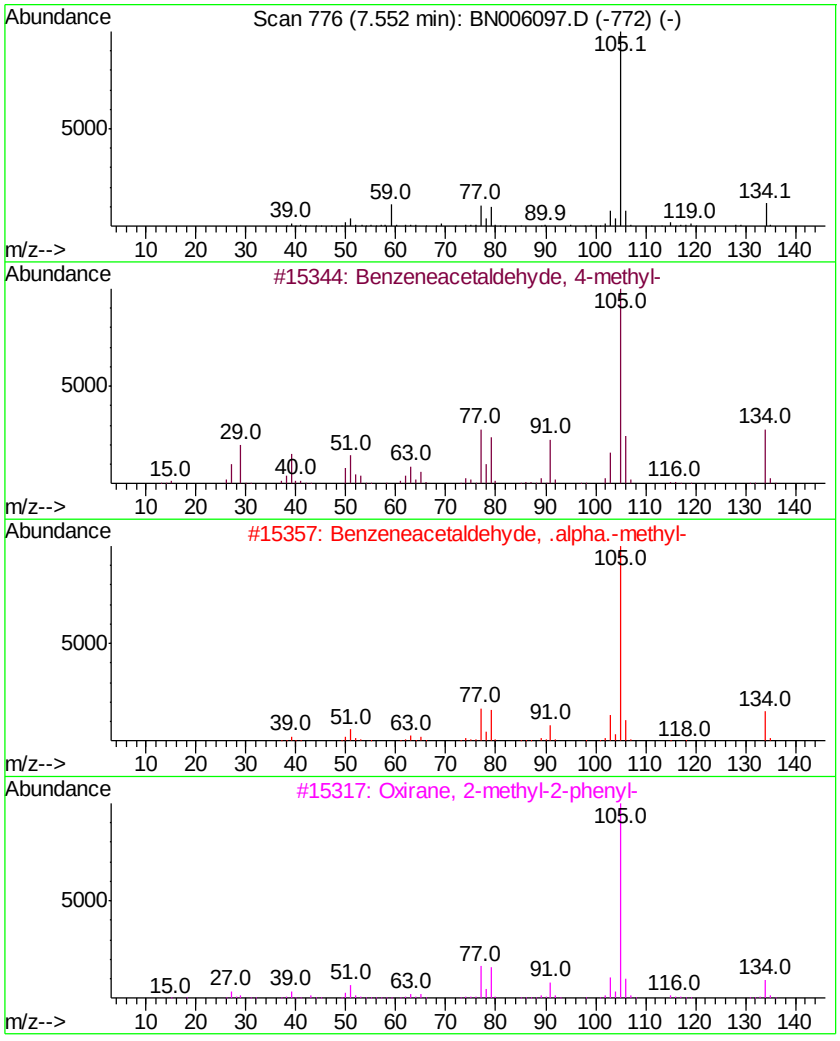
Quant Title : SVOA CALIBRATION

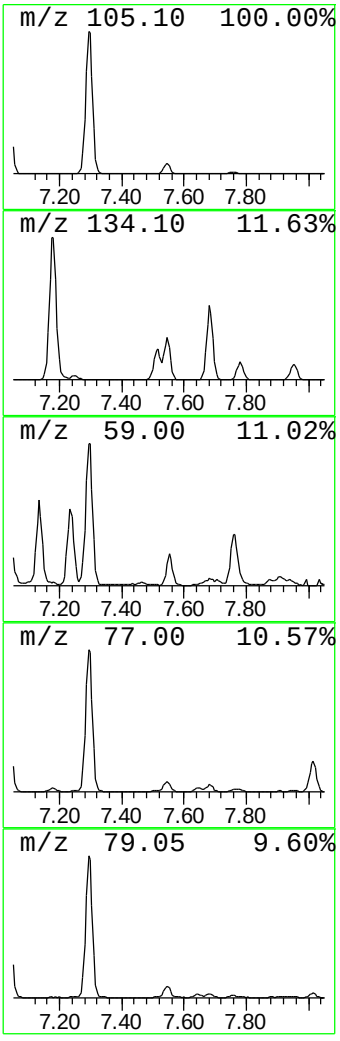
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzeneacetaldehyde, 4-methyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	5.92 ng/ul	566157	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetaldehyde, 4-methyl-	134	C9H10O	000104-09-6	80
2	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	74
3	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	74
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	72
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	64





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ALS Vial : 8 Sample Multiplier: 1

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Quant Title : SVOA CALIBRATION

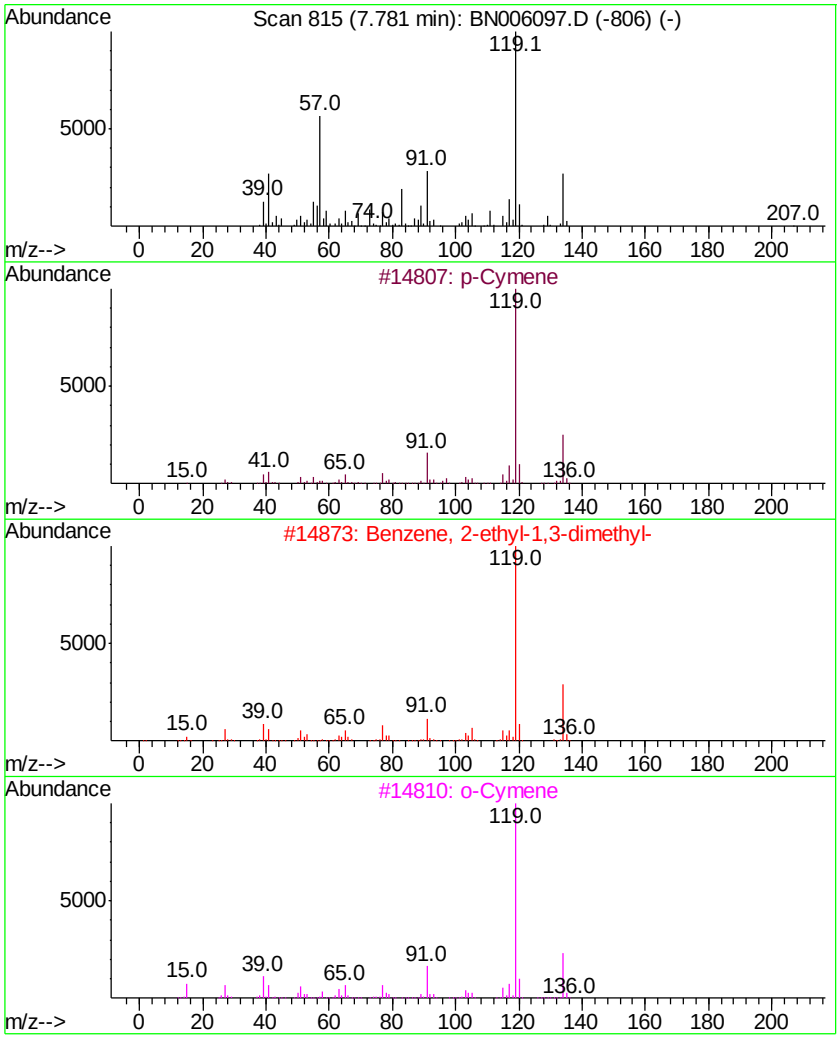
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

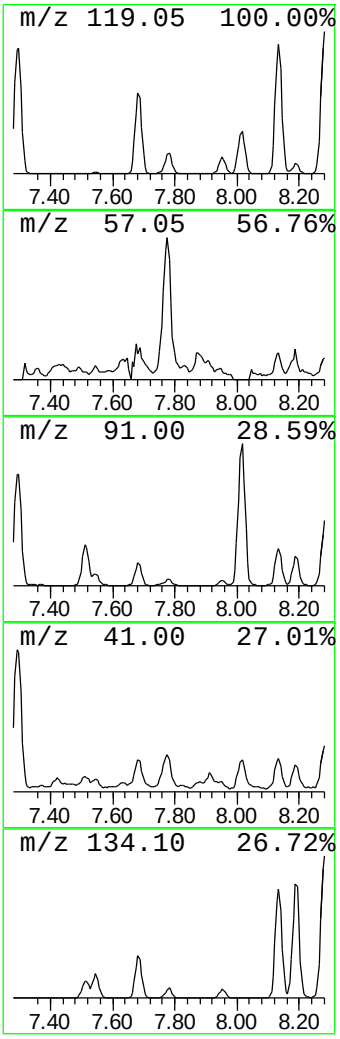
****

Peak Number 17 p-Cymene Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.78	6.37 ng/ul	609201	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	64
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
3	o-Cymene	134	C10H14	000527-84-4	64
4	p-Cymene	134	C10H14	000099-87-6	64
5	o-Cymene	134	C10H14	000527-84-4	64





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Quant Title : SVOA CALIBRATION

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

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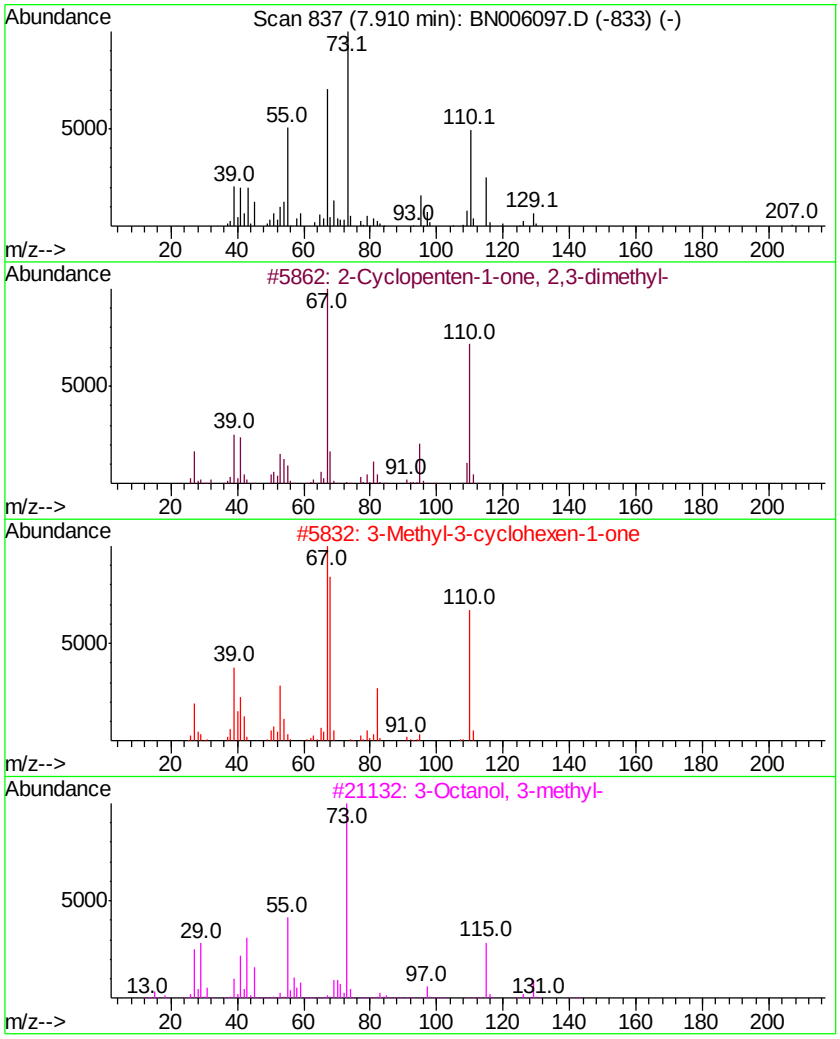
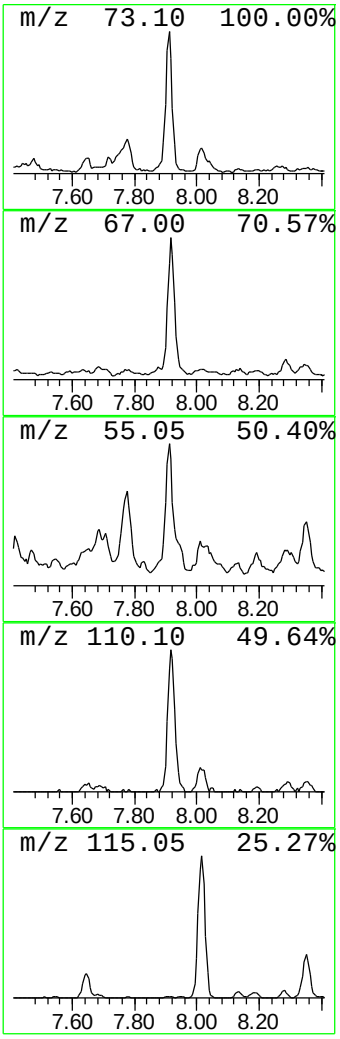
Peak Number 18 2-Cyclopenten-1-one, 2,3-di... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	3.28 ng/ul	313959	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclopenten-1-one, 2,3-dimethyl-	110	C7H10O	001121-05-7	53
2	3-Methyl-3-cyclohexen-1-one	110	C7H10O	031883-98-4	43
3	3-Octanol, 3-methyl-	144	C9H20O	005340-36-3	14
4	1,4-Pentadiene, 2-methyl-	82	C6H10	000763-30-4	11
5	Cyclopropane, 1,2-dimethyl-3-met...	82	C6H10	005070-00-8	11





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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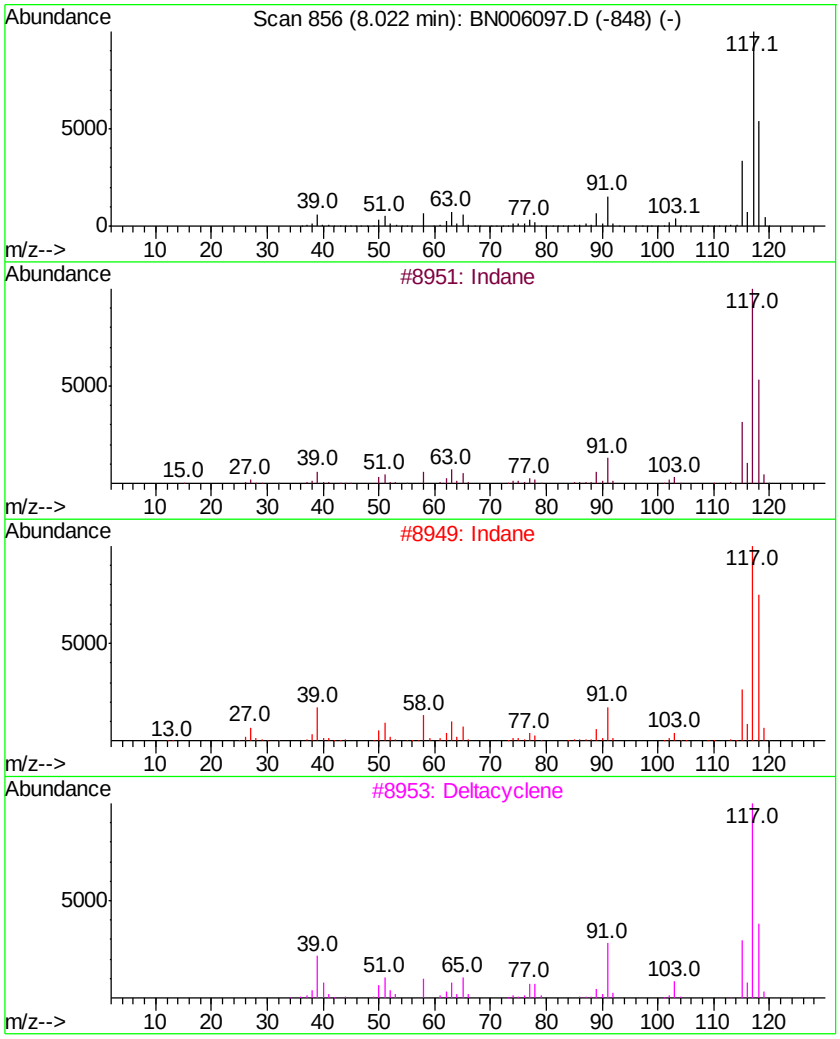
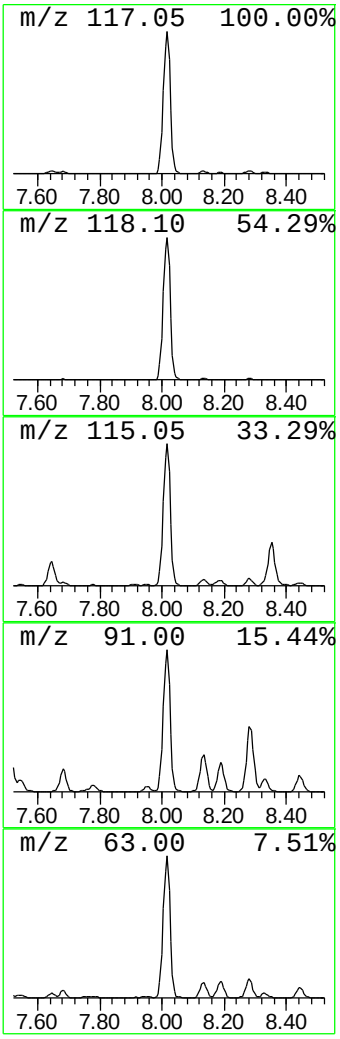
Peak Number 19 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	103.10 ng/ul	9857650	1,4-Dichlorobenzene-d4	7.65

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	Deltacyclene	118	C9H10	007785-10-6	72
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5	Indane	118	C9H10	000496-11-7	60





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Quant Title : SVOA CALIBRATION

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

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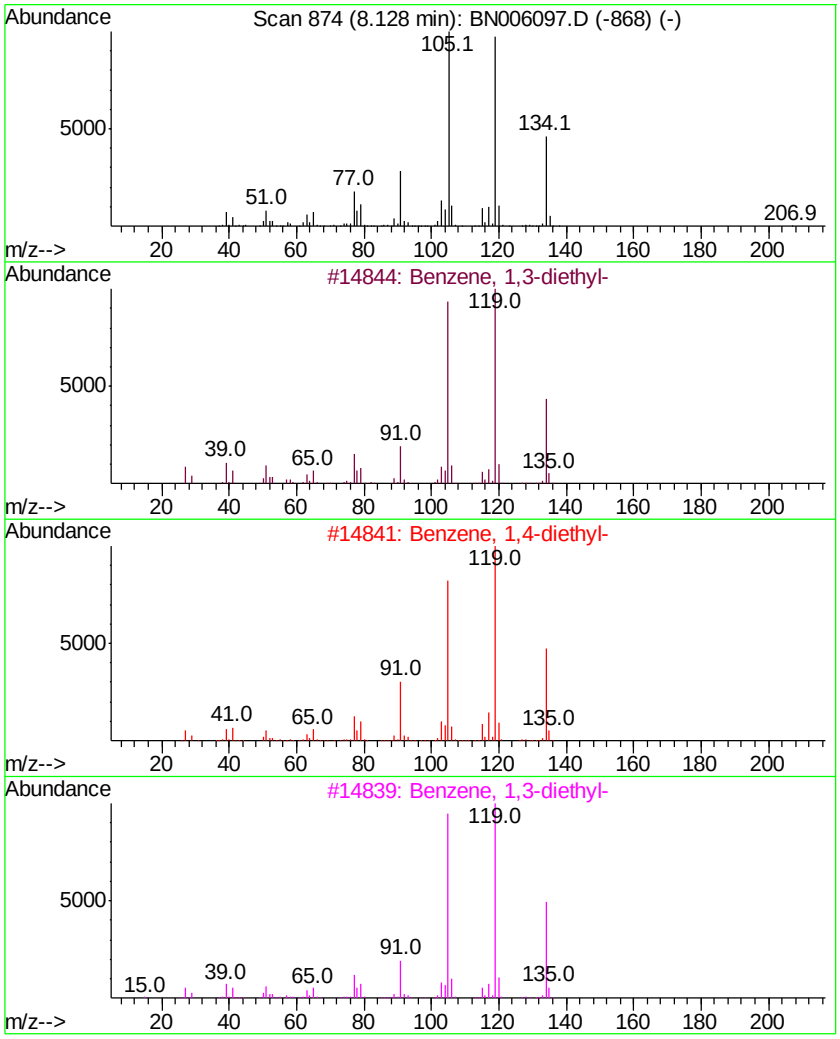
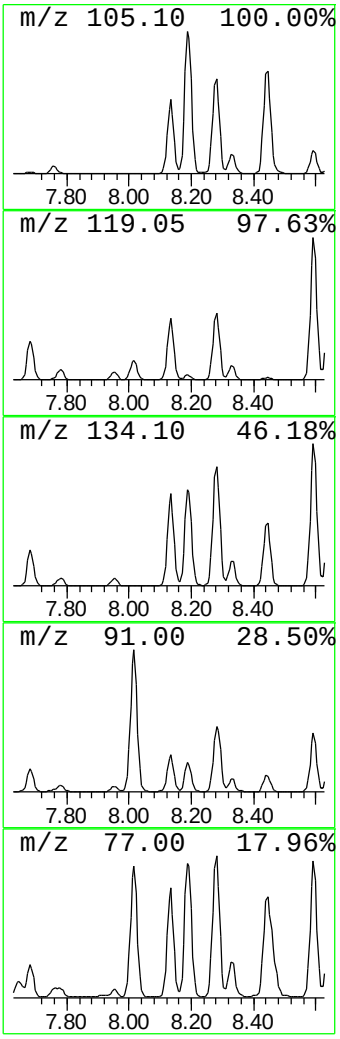
Peak Number 20 Benzene, 1,3-diethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.13	22.81 ng/ul	2181240	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95





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Quant Title : SVOA CALIBRATION

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

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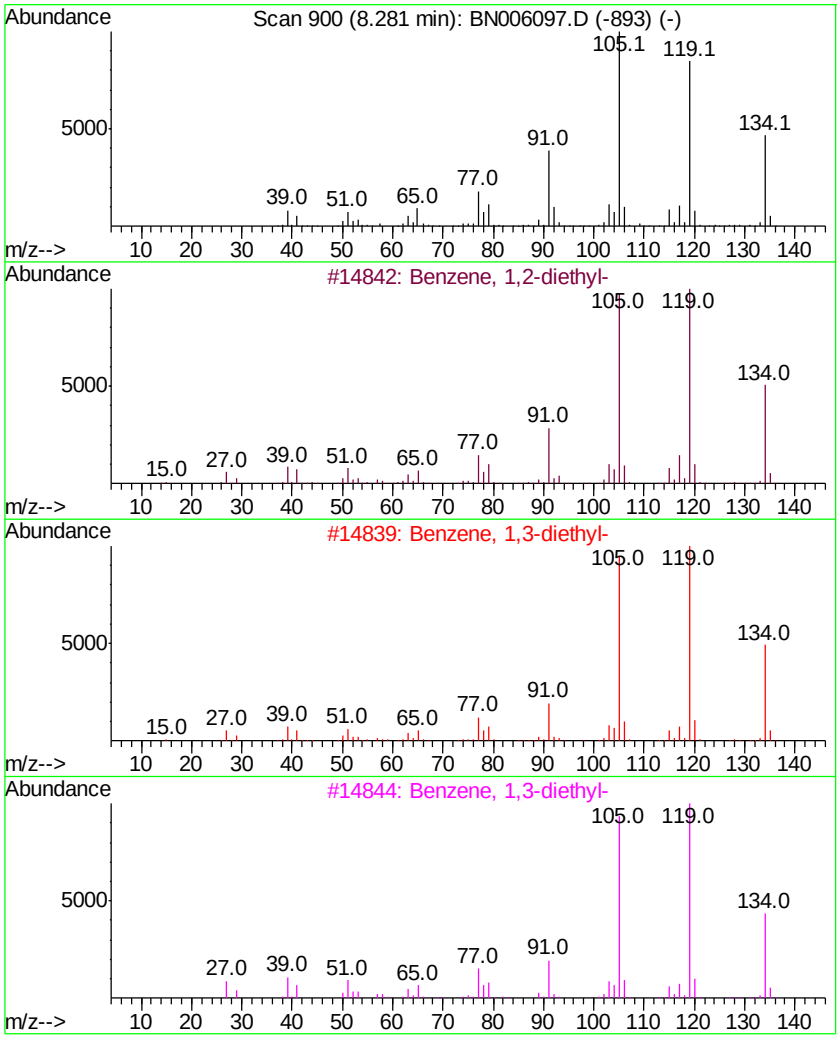
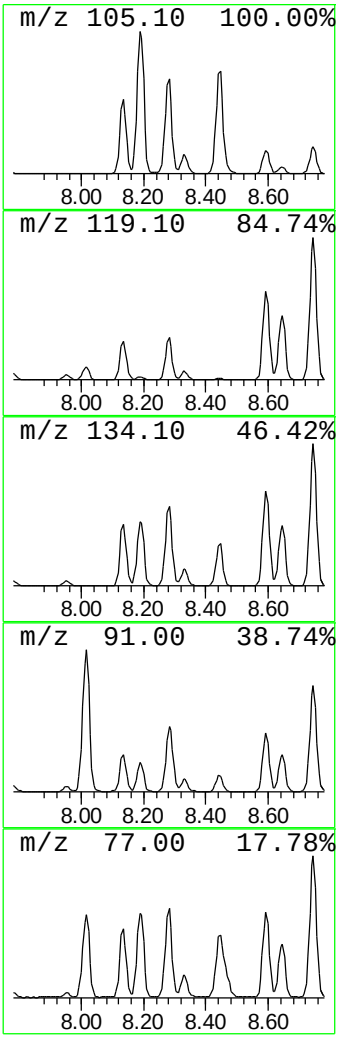
Peak Number 22 Benzene, 1,2-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.28	31.44 ng/ul	3005820	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93





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Instrument : BNA_N
ClientSampleId : C0C52

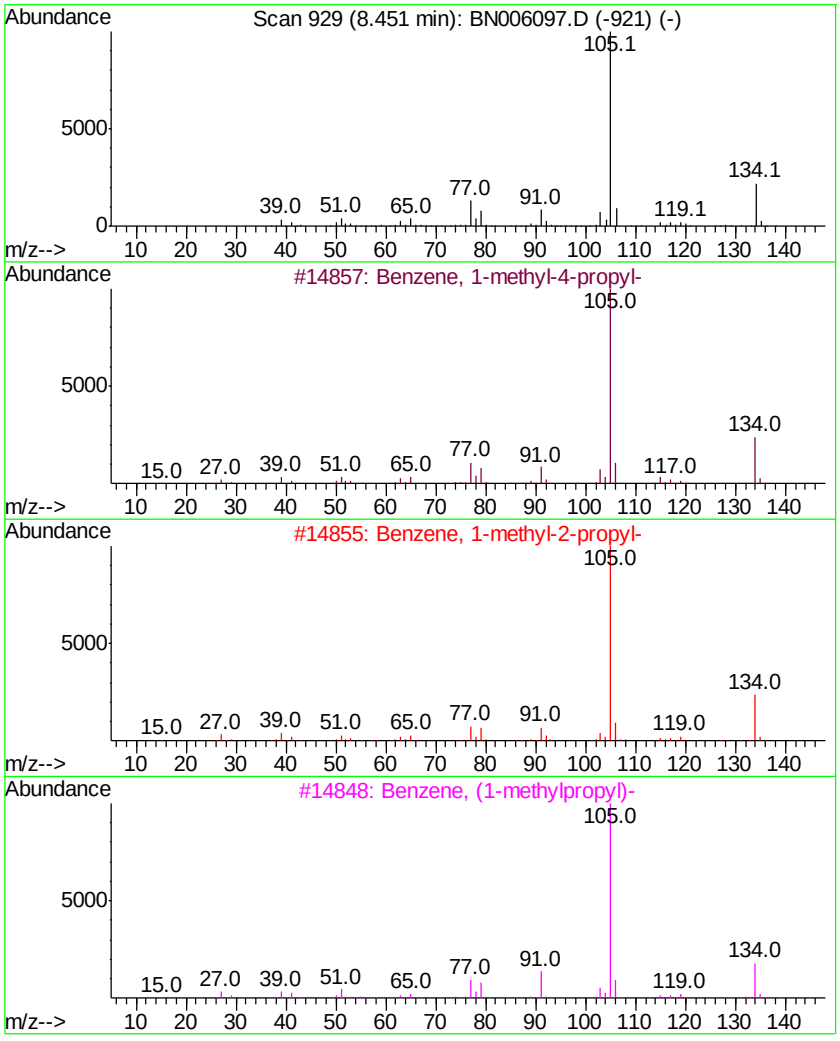
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

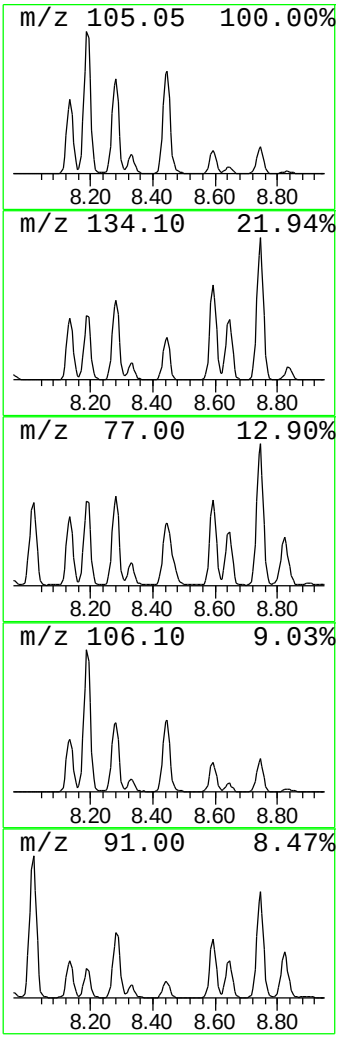
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1-methyl-4-propyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	17.61 ng/ul	1683940	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91





Library Search Compound Report

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Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

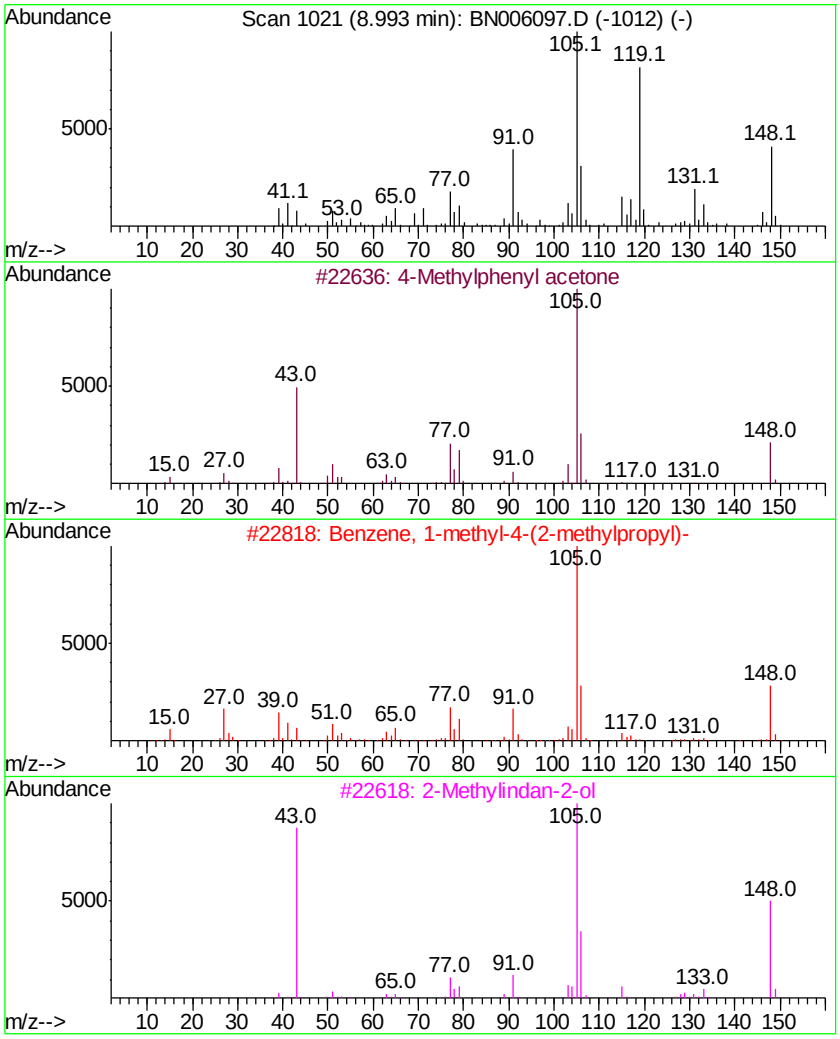
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

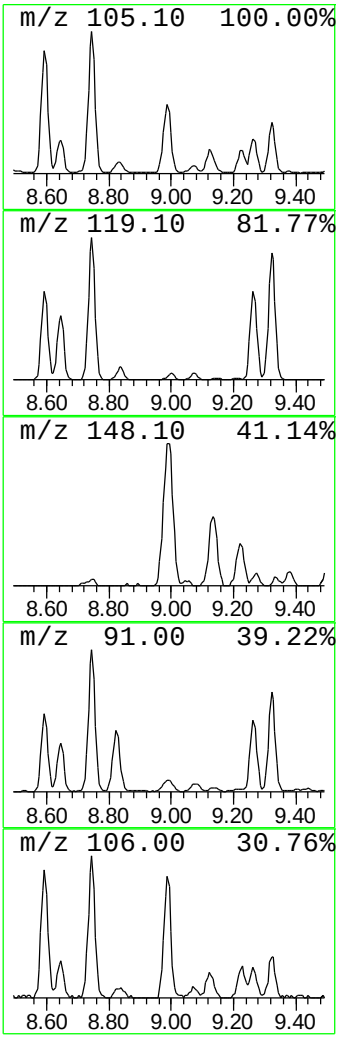
****

Peak Number 27 unknown-01 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.99	7.67 ng/ul	733464	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylphenyl acetone	148	C10H12O	002096-86-8	46
2	Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	41
3	2-Methylindan-2-ol	148	C10H12O	033223-84-6	41
4	Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	35
5	2-Butanone, 4-phenyl-	148	C10H12O	002550-26-7	30





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Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

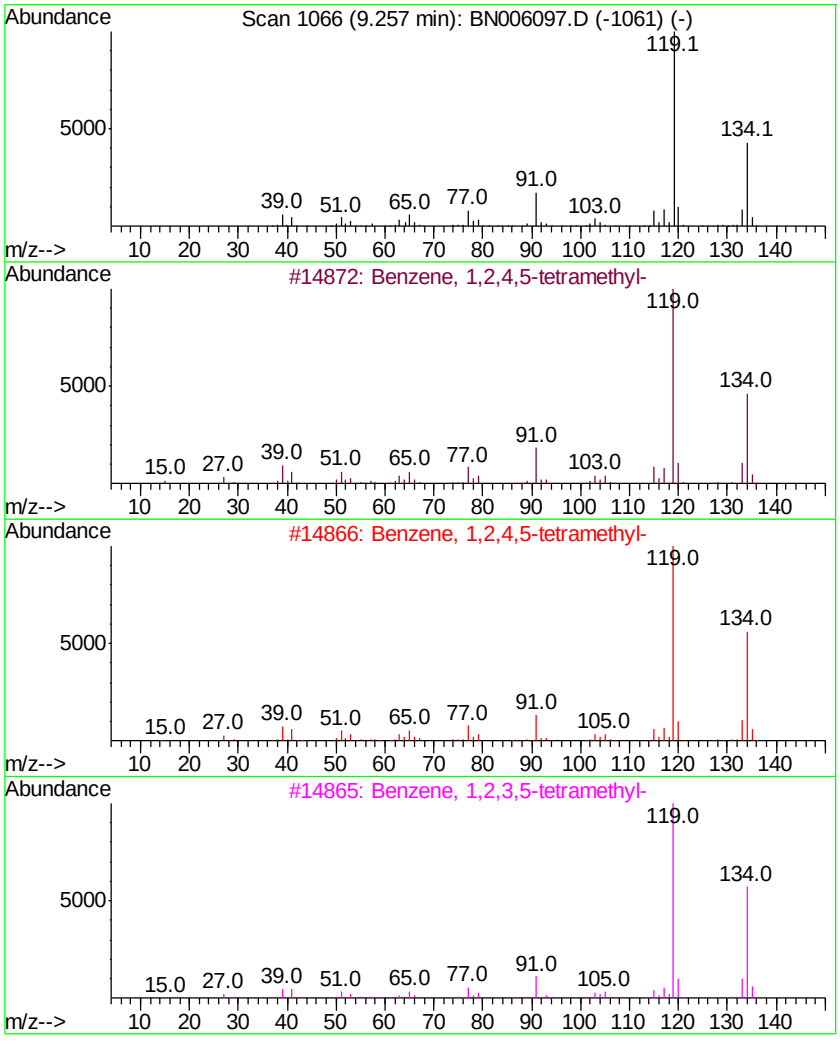
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

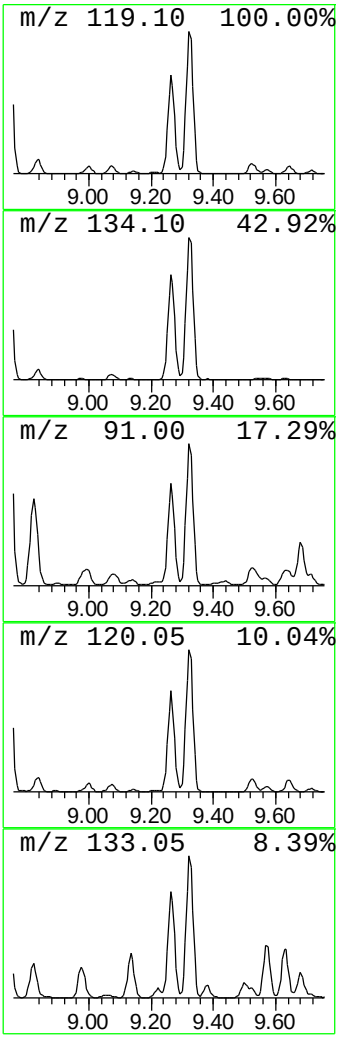
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.26	15.43 ng/ul	3031300	Naphthalene-d8	10.43

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	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
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





Library Search Compound Report

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Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

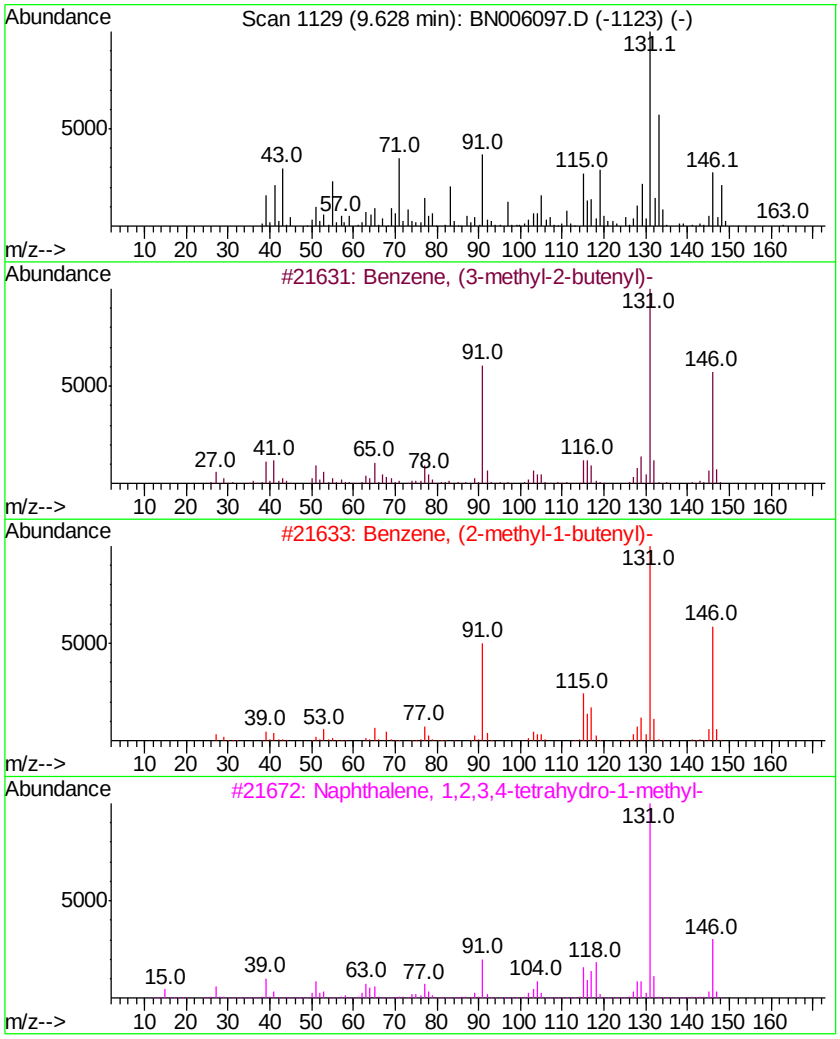
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

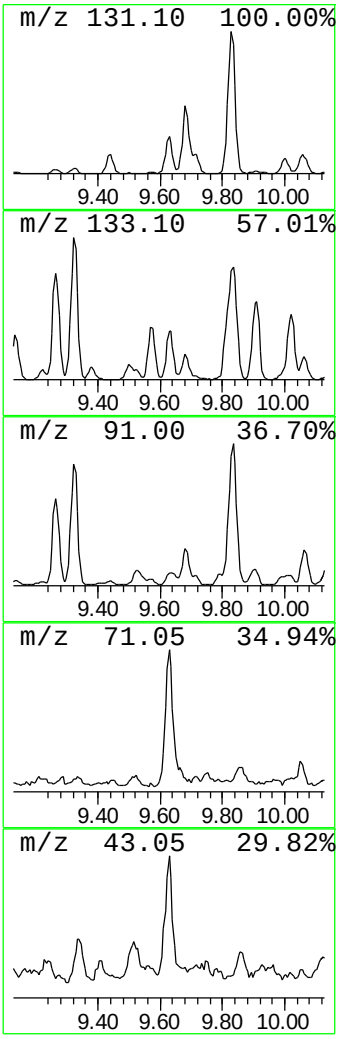
****

Peak Number 30 Benzene, (3-methyl-2-butenyl)- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	3.89 ng/ul	764509	Naphthalene-d8	10.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
2	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	90
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	78
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	78
5	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	60





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Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

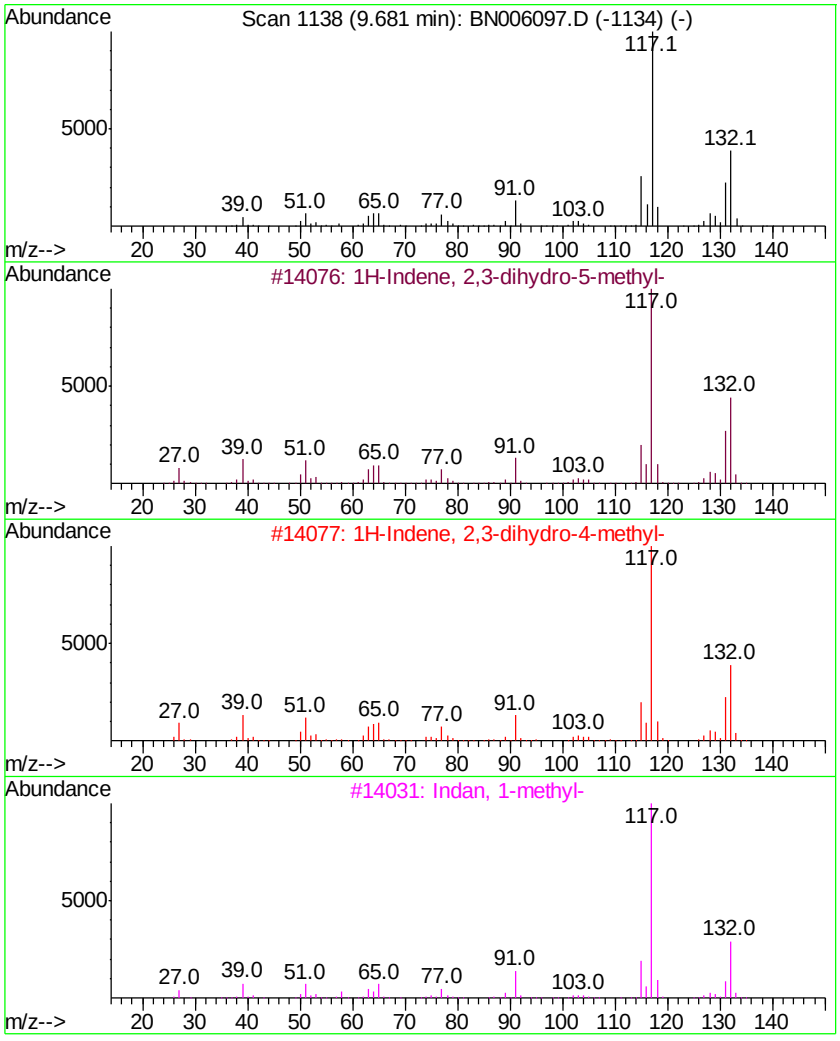
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

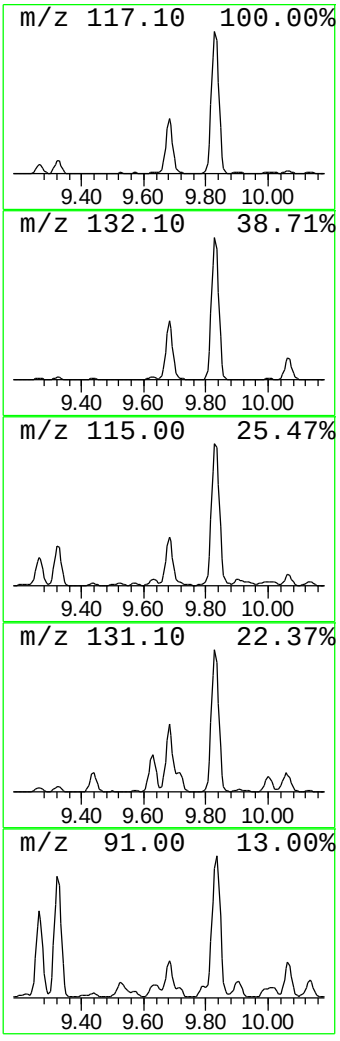
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	10.86 ng/ul	2134050	Naphthalene-d8	10.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81
5	Indan, 1-methyl-	132	C10H12	000767-58-8	80





Library Search Compound Report

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Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

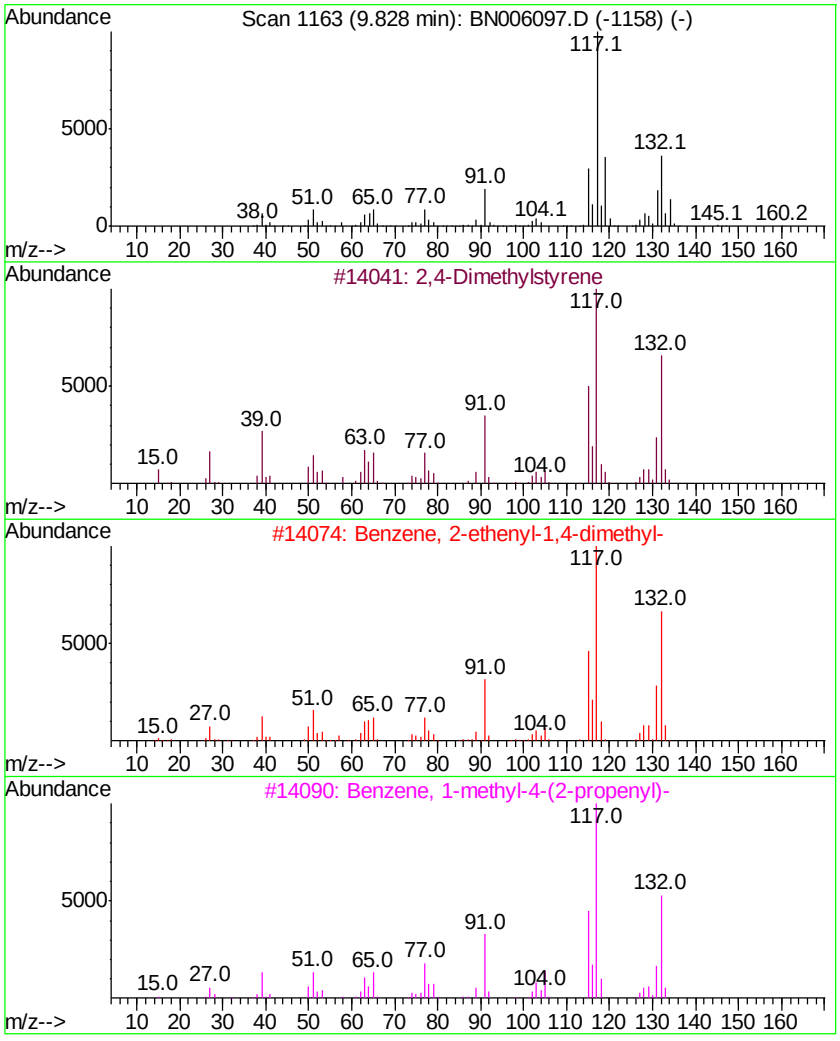
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

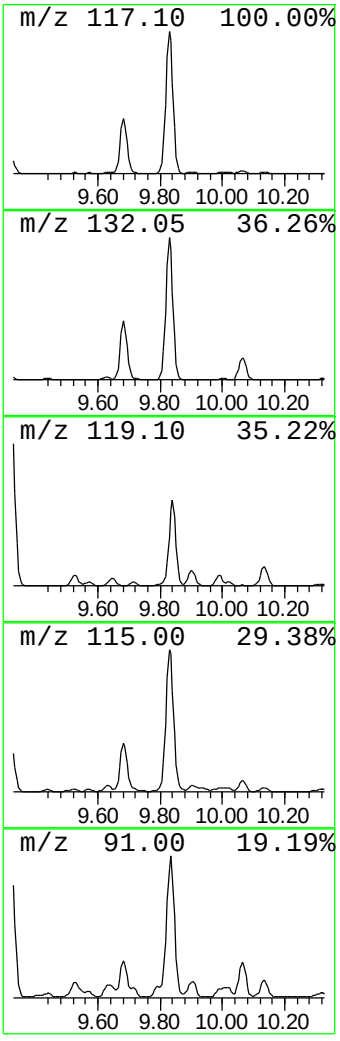
****

Peak Number 32 2,4-Dimethylstyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	37.94 ng/ul	7453210	Naphthalene-d8	10.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	87
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	86
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86





Library Search Compound Report

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Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

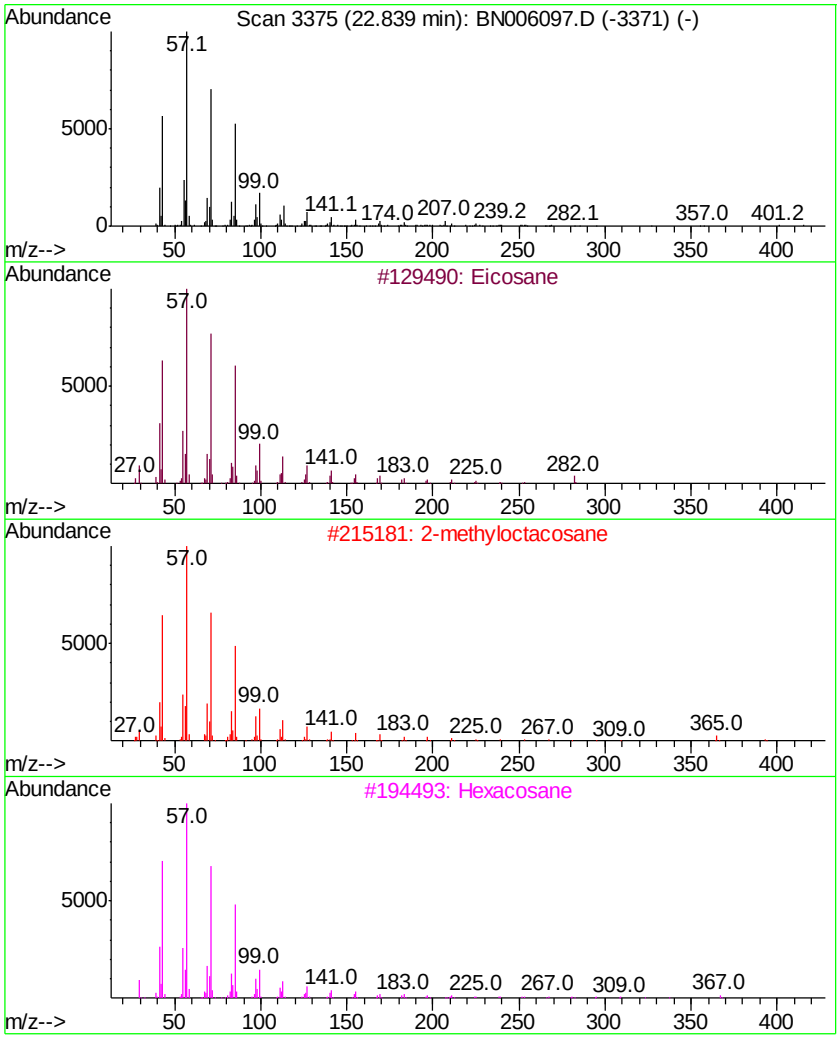
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

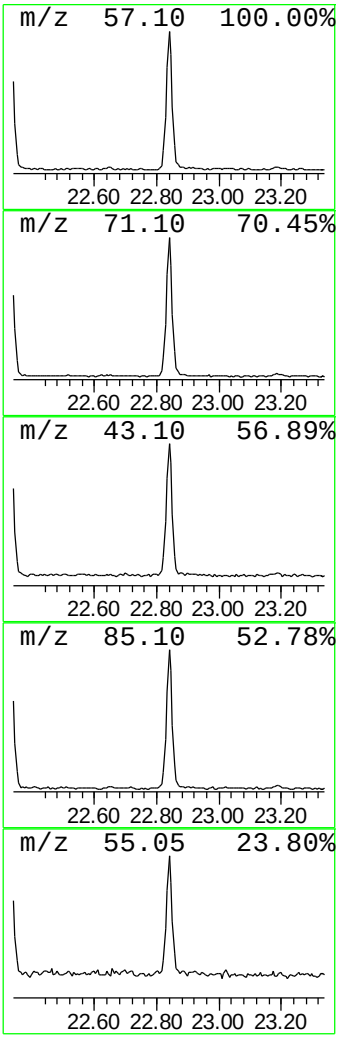
****

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.84	2.34 ng/ul	553907	Perylene-d12	23.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	91





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Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

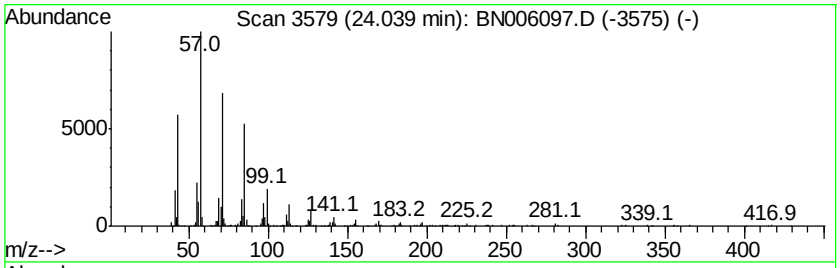
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

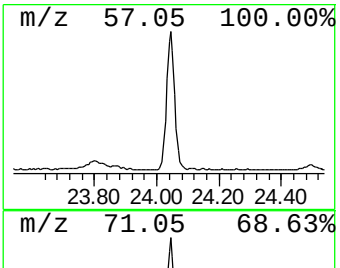
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

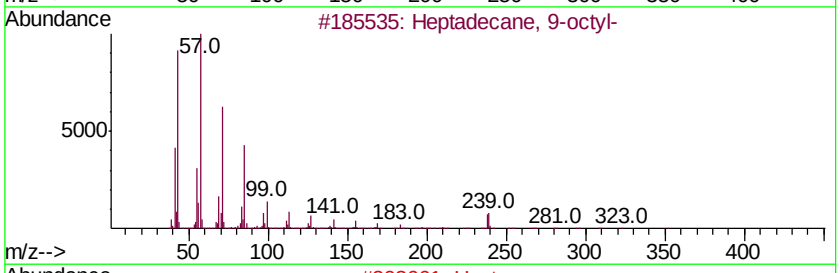
****
Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 48

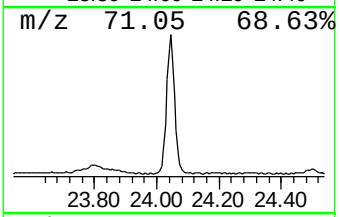
R.T.	EstConc	Area	Relative to ISTD	R.T.
24.04	3.32 ng/ul	788033	Perylene-d12	23.52

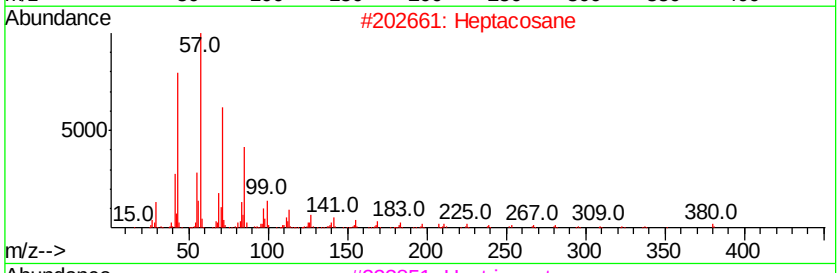
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2	Heptacosane	380	C27H56	000593-49-7	91
3	Hentriacontane	437	C31H64	000630-04-6	91
4	Octacosane	394	C28H58	000630-02-4	91
5	Tetratetracontane	619	C44H90	007098-22-8	91

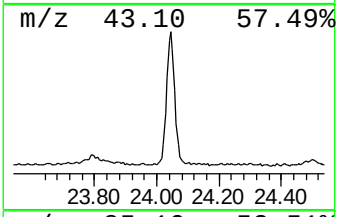


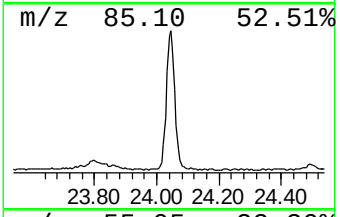


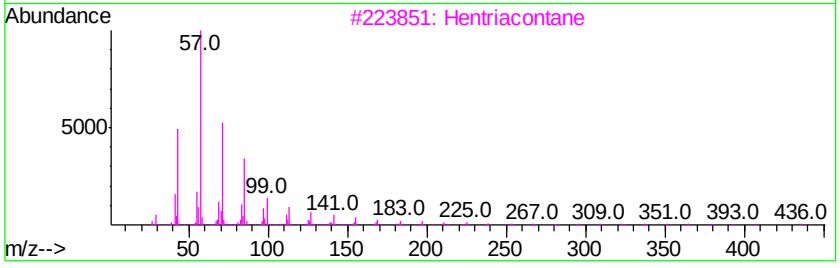


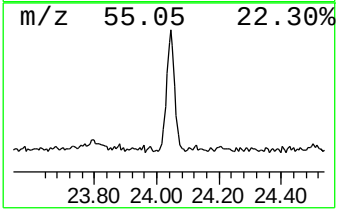












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Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

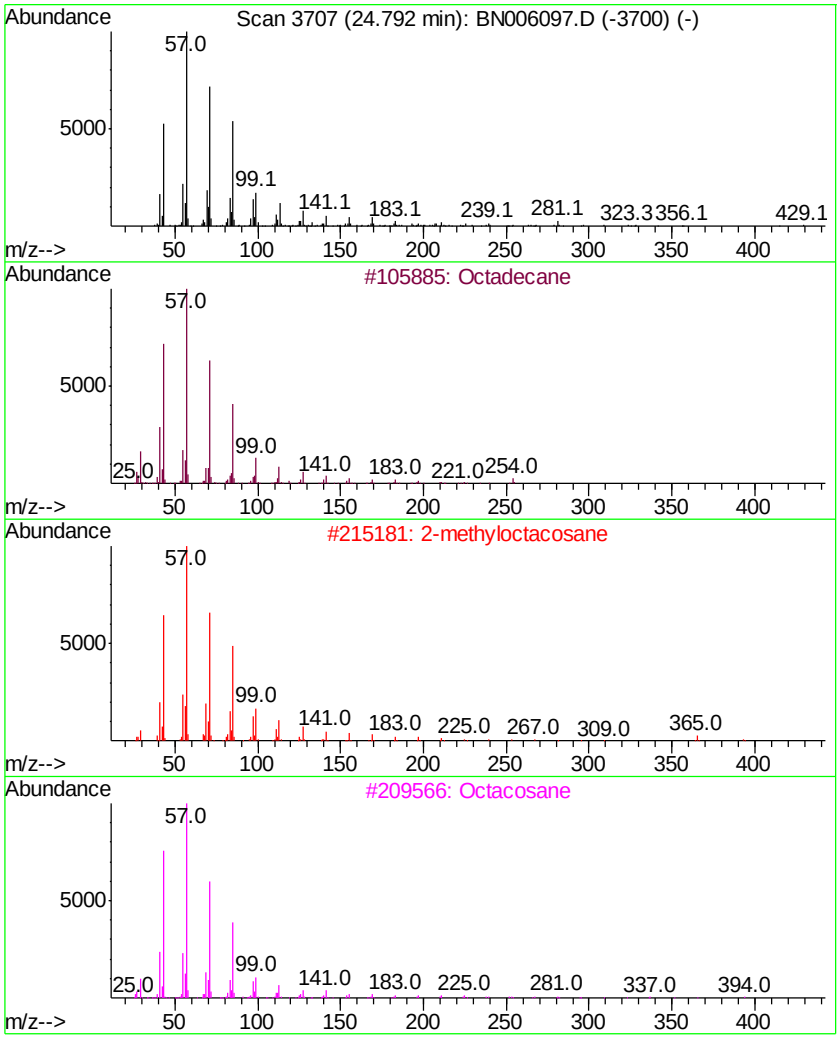
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

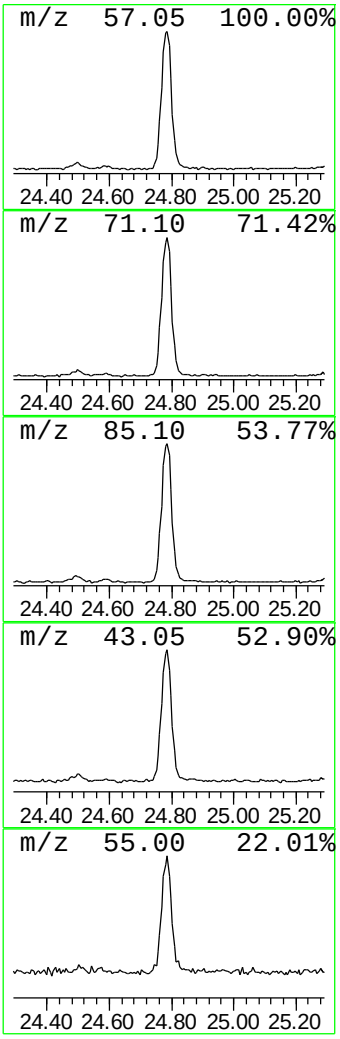
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.79	3.29 ng/ul	780790	Perylene-d12	23.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	93
2	2-methyloctacosane	408	C29H60	1000376-72-8	91
3	Octacosane	394	C28H58	000630-02-4	91
4	Heneicosane	296	C21H44	000629-94-7	90
5	Docosane, 11-decyl-	451	C32H66	055401-55-3	87





Library Search Compound Report

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Sample : K3045-08
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : C0C52

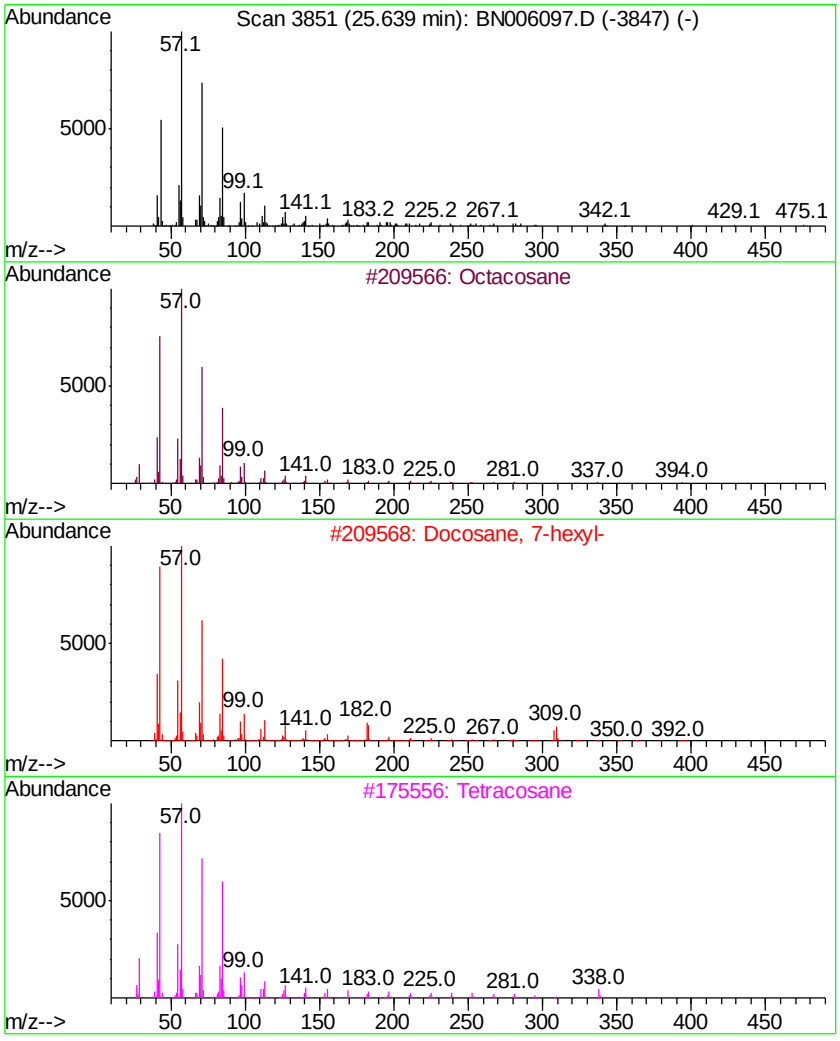
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060419MA.M
Quant Title : SVOA CALIBRATION

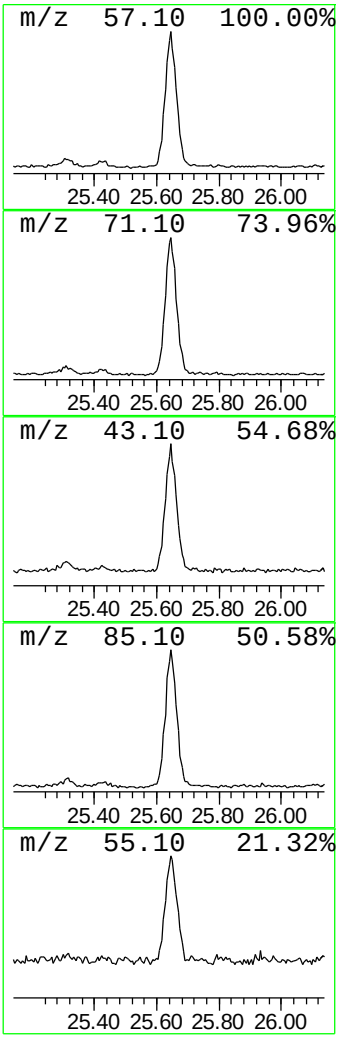
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.64	2.50 ng/ul	592330	Perylene-d12	23.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	90
2	Docosane, 7-hexyl-	394	C28H58	055373-86-9	87
3	Tetracosane	338	C24H50	000646-31-1	87
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	87
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87





Tentatively Identified Compound (LSC) summary

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Data File : BN006097.D
Acq On : 05 Jun 2019 13:52
Operator : JU/SJ
Sample : K3045-08
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ALS Vial : 8 Sample Multiplier: 1

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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
2-Pentanol, 2-met...	3.42	3.4	ng/ul	329194	1	7.65	1912340	20.0
(DEL) Alkane: Str...	3.73	3.9	ng/ul	369329	1	7.65	1912340	20.0
(DEL) Alkane: Cyc...	4.18	3.2	ng/ul	304626	1	7.65	1912340	20.0
3-Hexanol, 3-methyl-	4.83	4.9	ng/ul	470778	1	7.65	1912340	20.0
Benzene, (1-methy...	6.15	35.3	ng/ul	3376400	1	7.65	1912340	20.0
3,4-Dimethyl-3-he...	6.32	3.6	ng/ul	347924	1	7.65	1912340	20.0
Benzene, 1-ethyl-...	6.74	12.9	ng/ul	1234500	1	7.65	1912340	20.0
Benzene, 1,2,3-tr...	6.88	21.2	ng/ul	2028850	1	7.65	1912340	20.0
Benzene, 1-ethyl-...	7.04	57.4	ng/ul	5489990	1	7.65	1912340	20.0
Benzene, (2-methy...	7.51	4.2	ng/ul	398995	1	7.65	1912340	20.0
Benzeneacetaldehy...	7.55	5.9	ng/ul	566157	1	7.65	1912340	20.0
p-Cymene	7.78	6.4	ng/ul	609201	1	7.65	1912340	20.0
2-Cyclopenten-1-o...	7.91	3.3	ng/ul	313959	1	7.65	1912340	20.0
Indane	8.02	103.1	ng/ul	9857650	1	7.65	1912340	20.0
Benzene, 1,3-diet...	8.13	22.8	ng/ul	2181240	1	7.65	1912340	20.0
Benzene, 1,2-diet...	8.28	31.4	ng/ul	3005820	1	7.65	1912340	20.0
Benzene, 1-methyl...	8.45	17.6	ng/ul	1683940	1	7.65	1912340	20.0
unknown-01	8.99	7.7	ng/ul	733464	1	7.65	1912340	20.0
Benzene, 1,2,4,5-...	9.26	15.4	ng/ul	3031300	2	10.43	3928740	20.0
Benzene, (3-methy...	9.63	3.9	ng/ul	764509	2	10.43	3928740	20.0
1H-Indene, 2,3-di...	9.68	10.9	ng/ul	2134050	2	10.43	3928740	20.0
2,4-Dimethylstyrene	9.83	37.9	ng/ul	7453210	2	10.43	3928740	20.0
(DEL) Alkane: Str...	22.84	2.3	ng/ul	553907	6	23.52	4743400	20.0
(DEL) Alkane: Str...	24.04	3.3	ng/ul	788033	6	23.52	4743400	20.0
(DEL) Alkane: Str...	24.79	3.3	ng/ul	780790	6	23.52	4743400	20.0
(DEL) Alkane: Str...	25.64	2.5	ng/ul	592330	6	23.52	4743400	20.0