LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056831.D
Acq On : 4 Mar 2023 10:24
Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBAD4

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_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

Signal : TIC: BG056831.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.137	99	102	114	rBV	55214	97019	13.93%	1.251%
2	7.080	599	603	608	rBV4	8262	13493	1.94%	0.174%
3	7.538	675	681	692	rVB	96071	170010	24.42%	2.192%
4	7.721	706	712	722	rBV	66715	109083	15.67%	1.407%
5	7.938	742	749	756	rVB	85276	147821	21.23%	1.906%
6	8.420	824	831	837	rVB	56917	92584	13.30%	1.194%
7	8.725	878	883	888	rVB3	15195	22950	3.30%	0.296%
8	9.107	941	948	955	rBV	120994	210005	30.16%	2.708%
9	9.166	955	958	964	rVB4	2779	4198	0.60%	0.054%
10	9.589	1023	1030	1039	rVB	99540	172677	24.80%	2.227%
11	9.965	1092	1094	1099	rVB4	1430	1613	0.23%	0.021%
12	10.318	1147	1154	1162	rBV2	86771	151422	21.75%	1.953%
13	10.664	1211	1213	1216	rVB2	1289	969	0.14%	0.012%
14	10.858	1239	1246	1256	rBV2	128012	225893	32.44%	2.913%
15	11.017	1269	1273	1279	rVB3	2108	3308	0.48%	0.043%
16	11.252	1307	1313	1322	rVB	80391	146937	21.10%	1.895%
17	11.375	1327	1334	1343	rVB	109295	191992	27.57%	2.476%
18	12.039	1444	1447	1452	rVB3	796	968	0.14%	0.012%
19	12.715	1558	1562	1566	rBV2	484	981	0.14%	0.013%
20	12.803	1574	1577	1582	rVB3	1965	2638	0.38%	0.034%
21	12.979	1602	1607	1614	rVB2	6379	9604	1.38%	0.124%
22	13.390	1673	1677	1680	rVB3	1521	1924	0.28%	0.025%
23	13.813	1745	1749	1754	rVB3	1979	3054	0.44%	0.039%
24	13.896	1759	1763	1767	rBV3	8959	12555	1.80%	0.162%
25	14.178	1807	1811	1818	rVB2	17494	24392	3.50%	0.315%
26	14.348	1834	1840	1845	rBV	171186	259367	37.25%	3.345%
27	14.407	1845	1850	1857	rVB	246353	355302	51.03%	4.582%
28	14.718	1896	1903	1911	rVB	277646	430007	61.76%	5.545%
29	14.848	1922	1925	1928	rBV2	1450	1303	0.19%	0.017%
30	14.871	1928	1929	1937	rVB3	1489	1244	0.18%	0.016%
31	14.965	1943	1945	1948	rBV2	2093	1852	0.27%	0.024%
32	15.018	1948	1954	1960	rVV	149194	230275	33.07%	2.969%
33	15.071	1960	1963	1968	rVB4	4572	5256	0.75%	0.068%
34	15.177	1974	1981	1987	rBV	120267	189036	27.15%	2.438%
35	15.247	1990	1993	1996	rVV3	4258	4445	0.64%	0.057%
36	15.723	2071	2074	2077	rBV3	1409	1919	0.28%	0.025%

LSC Area Percent Report

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Sample : 01711-20
Misc :
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Instrument :
BNA_G
ClientSampleId :
DBAD4

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_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

37	15.887	2099	2102	2106	rBV2	1256	1551	0.22%	0.020%
38	16.005	2115	2122	2129	rVB	347925	539136	77.43%	6.952%
39	16.099	2131	2138	2144	rBV	123319	182694	26.24%	2.356%
40	16.246	2159	2163	2166	rVB2	1010	1678	0.24%	0.022%
41	16.352	2174	2181	2188	rVB	23754	35403	5.08%	0.457%
42	16.410	2188	2191	2193	rBV2	1576	1811	0.26%	0.023%
43	16.463	2196	2200	2204	rVB2	10175	14293	2.05%	0.184%
44	16.557	2213	2216	2220	rVV	927	1146	0.16%	0.015%
45	16.622	2223	2227	2230	rVB2	2306	3312	0.48%	0.043%
46	16.716	2238	2243	2245	rVB2	1166	1540	0.22%	0.020%
47	17.356	2348	2352	2354	rVB2	824	1092	0.16%	0.014%
48	17.450	2367	2368	2372	rVB2	1339	1244	0.18%	0.016%
49	17.591	2390	2392	2395	rBV	1382	1327	0.19%	0.017%
50	17.674	2401	2406	2407	rBV2	771	957	0.14%	0.012%
51	17.697	2407	2410	2412	rVV2	1501	1651	0.24%	0.021%
52	17.750	2413	2419	2427	rVB	201899	311137	44.69%	4.012%
53	17.850	2429	2436	2443	rBV2	380009	581189	83.47%	7.495%
54	17.915	2443	2447	2448	rBV	2245	2908	0.42%	0.037%
55	17.938	2448	2451	2454	rVB4	5174	6732	0.97%	0.087%
56	17.997	2460	2461	2466	rVB2	1189	974	0.14%	0.013%
57	18.308	2510	2514	2517	rVB2	738	1022	0.15%	0.013%
58	18.461	2538	2540	2541	rBV	1710	1606	0.23%	0.021%
59	18.490	2541	2545	2551	rVV4	9469	15139	2.17%	0.195%
60	18.590	2558	2562	2575	rBV	49809	89859	12.91%	1.159%
61	18.808	2594	2599	2604	rBV2	42305	57834	8.31%	0.746%
62	19.019	2630	2635	2641	rVB5	10304	19020	2.73%	0.245%
63	19.148	2654	2657	2659	rVB2	2414	1985	0.29%	0.026%
64	19.201	2661	2666	2670	rVV2	21925	28656	4.12%	0.370%
65	19.331	2687	2688	2693	rVB2	1725	1859	0.27%	0.024%
66	19.677	2742	2747	2750	rBV4	5095	7503	1.08%	0.097%
67	19.748	2754	2759	2763	rVB3	5462	10143	1.46%	0.131%
68	19.783	2763	2765	2767	rBV	1376	1315	0.19%	0.017%
69	19.818	2768	2771	2772	rBV	1682	1595	0.23%	0.021%
70	19.842	2772	2775	2778	rVB2	6539	7380	1.06%	0.095%
71	19.977	2797	2798	2800	rBV	1362	1133	0.16%	0.015%
72	20.059	2810	2812	2815	rBV3	1449	1231	0.18%	0.016%
73	20.112	2815	2821	2827	rBV	499949	696287	100.00%	8.979%
74	20.212	2835	2838	2843	rBV	17127	23827	3.42%	0.307%
75	20.435	2872	2876	2879	rBV4	3570	4708	0.68%	0.061%
76	20.529	2889	2892	2896	rVB2	7165	8390	1.20%	0.108%

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Instrument :
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ClientSampleId :
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Integration Parameters: LSCINT.P
Integrator: RTE
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Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

77	20.564	2896	2898	2899	rBV2	3172	1690	0.24%	0.022%
78	20.611	2903	2906	2911	rVB2	20044	20719	2.98%	0.267%
79	20.664	2913	2915	2919	rBV4	1486	1388	0.20%	0.018%
80	20.805	2937	2939	2945	rVB5	4610	4705	0.68%	0.061%
81	20.876	2947	2951	2958	rBV6	11209	22242	3.19%	0.287%
82	20.935	2958	2961	2966	rBV	23993	30868	4.43%	0.398%
83	21.058	2979	2982	2986	rBV	6383	7448	1.07%	0.096%
84	21.281	3018	3020	3023	rVB4	3617	3211	0.46%	0.041%
85	21.381	3035	3037	3039	rBV3	2146	1421	0.20%	0.018%
86	21.457	3045	3050	3058	rVB	147519	208273	29.91%	2.686%
87	21.510	3058	3059	3062	rBV3	2440	2639	0.38%	0.034%
88	21.534	3062	3063	3066	rVB2	2823	2724	0.39%	0.035%
89	21.634	3075	3080	3091	rBV2	60620	109003	15.65%	1.406%
90	22.063	3146	3153	3160	rVB	193622	339626	48.78%	4.380%
91	22.221	3177	3180	3188	rVB6	5839	9164	1.32%	0.118%
92	23.520	3399	3401	3402	rBV2	2438	1707	0.25%	0.022%
93	23.849	3454	3457	3462	rVB4	4331	7347	1.06%	0.095%
94	25.347	3700	3712	3723	rVB	230552	672624	96.60%	8.674%
95	25.588	3743	3753	3763	rVB2	112786	336244	48.29%	4.336%
96	26.040	3828	3830	3831	rBV2	1869	1380	0.20%	0.018%

Sum of corrected areas: 7754786

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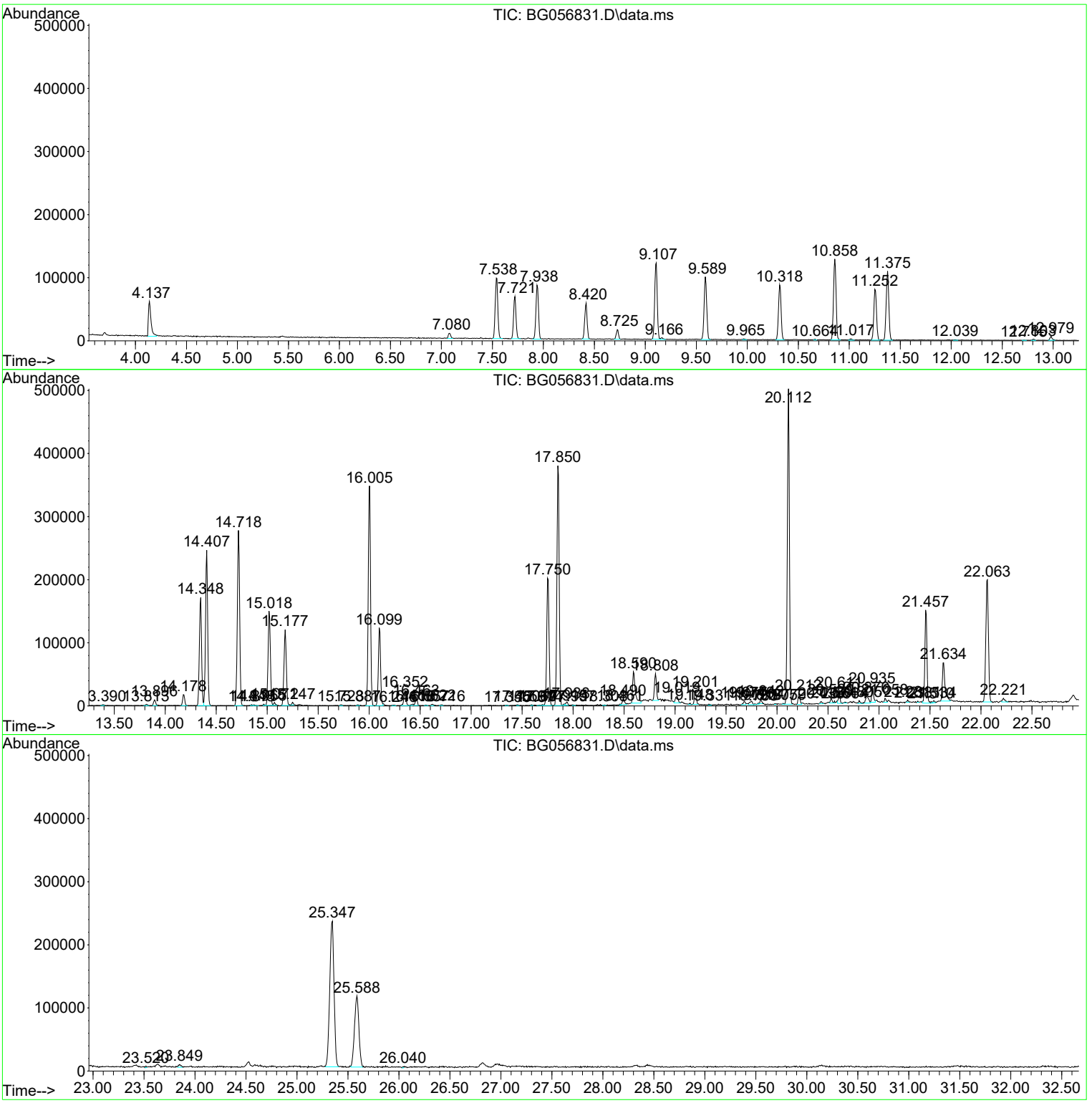
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
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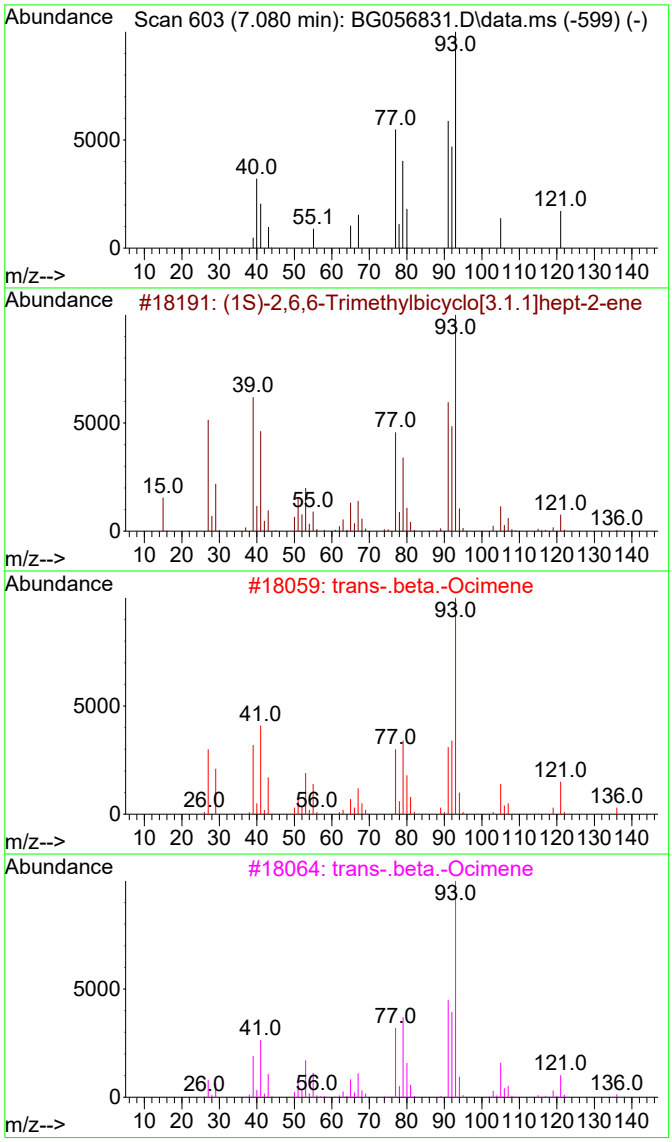
Quant Title : SVOA CALIBRATION

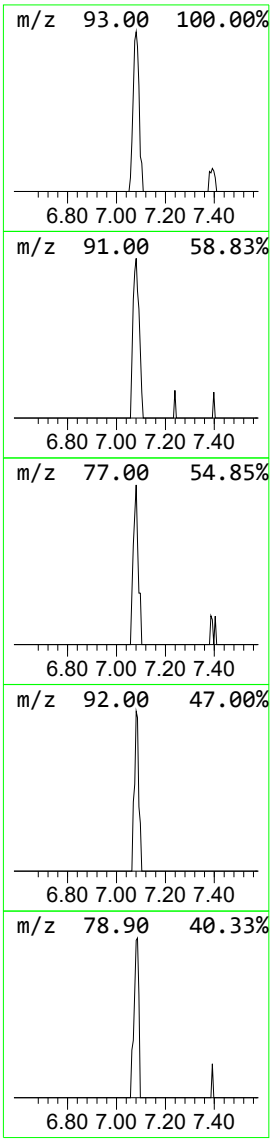
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (1S)-2,6,6-Trimethylbicyclo... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.080	2.91 ng/ul	13493	1,4-Dichlorobenzene-d4	8.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	83
2	trans-.beta.-Ocimene	136	C10H16	003779-61-1	72
3	trans-.beta.-Ocimene	136	C10H16	003779-61-1	72
4	1,3,6-Heptatriene, 5-methyl-, (E)-	108	C8H12	000818-48-4	59
5	trans-.beta.-Ocimene	136	C10H16	003779-61-1	45





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ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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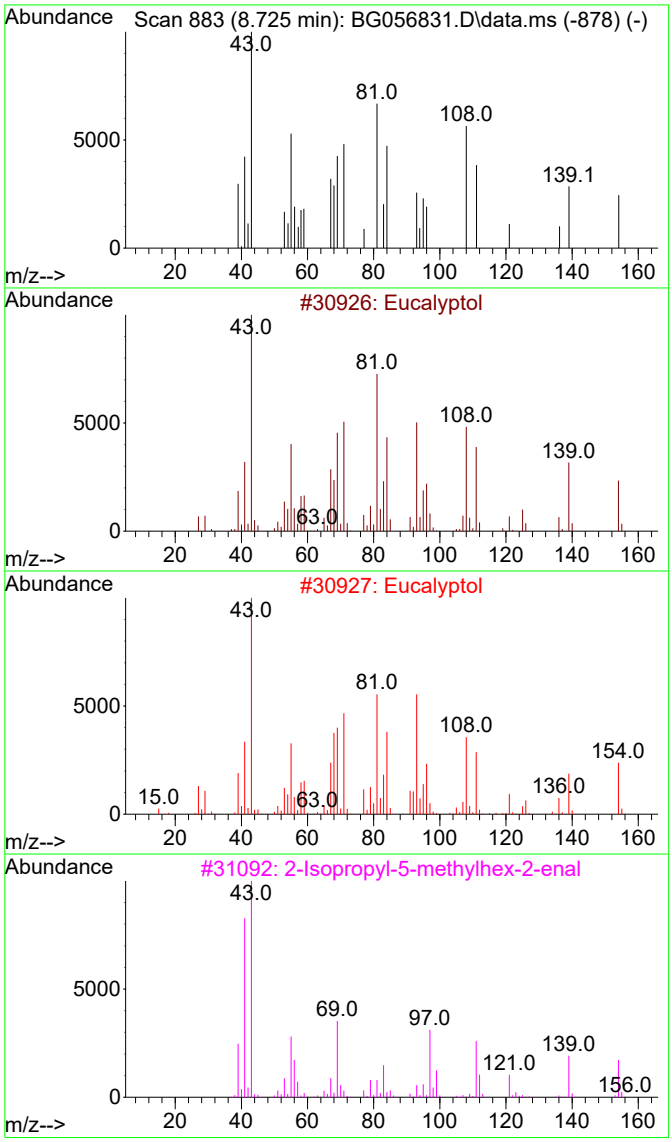
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

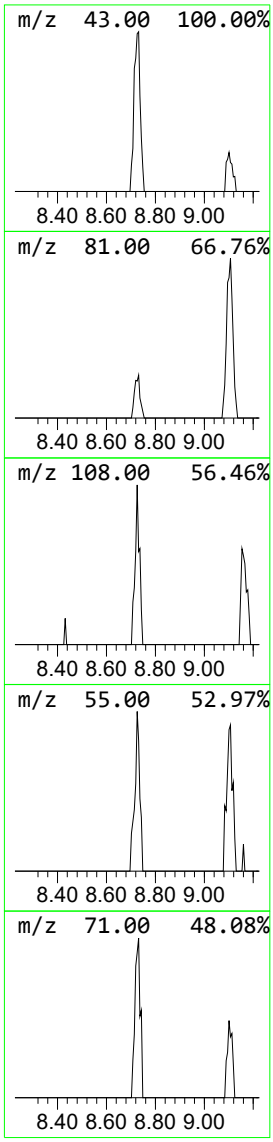
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Eucalyptol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.725	4.96 ng/ul	22950	1,4-Dichlorobenzene-d4	8.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	94
2	Eucalyptol	154	C10H18O	000470-82-6	58
3	2-Isopropyl-5-methylhex-2-enal	154	C10H18O	035158-25-9	25
4	5-Hepten-2-one, 6-methyl-	126	C8H14O	000110-93-0	22
5	4-Isopropyl-1-methylcyclohex-2-enol	154	C10H18O	000619-62-5	17





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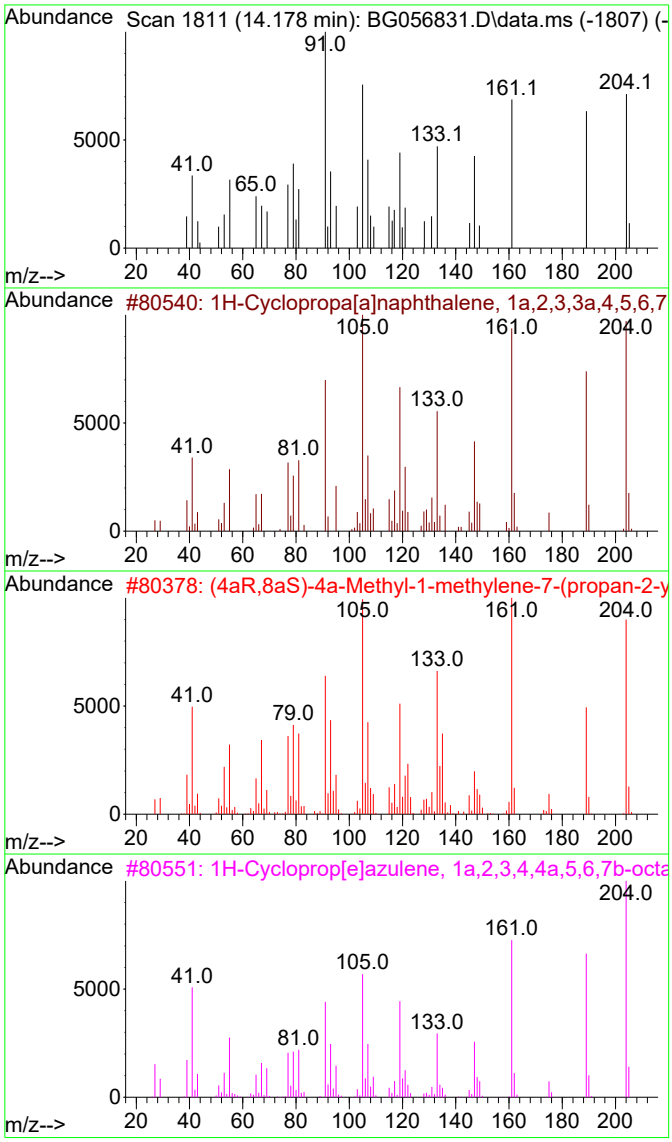
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

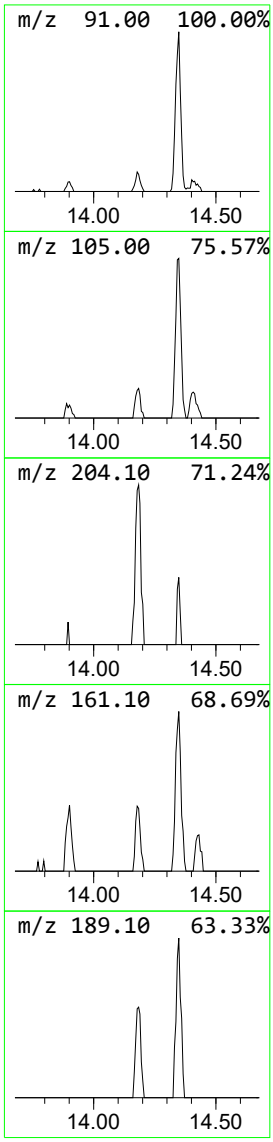
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1H-Cyclopropa[a]naphthalene... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.178	2.12 ng/ul	24392	Acenaphthene-d10	15.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	000489-29-2	96
2	(4aR,8aS)-4a-Methyl-1-methylene-...	204	C15H24	058893-88-2	93
3	1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	92
4	.beta.-Guaiene	204	C15H24	000088-84-6	91
5	1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	91





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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

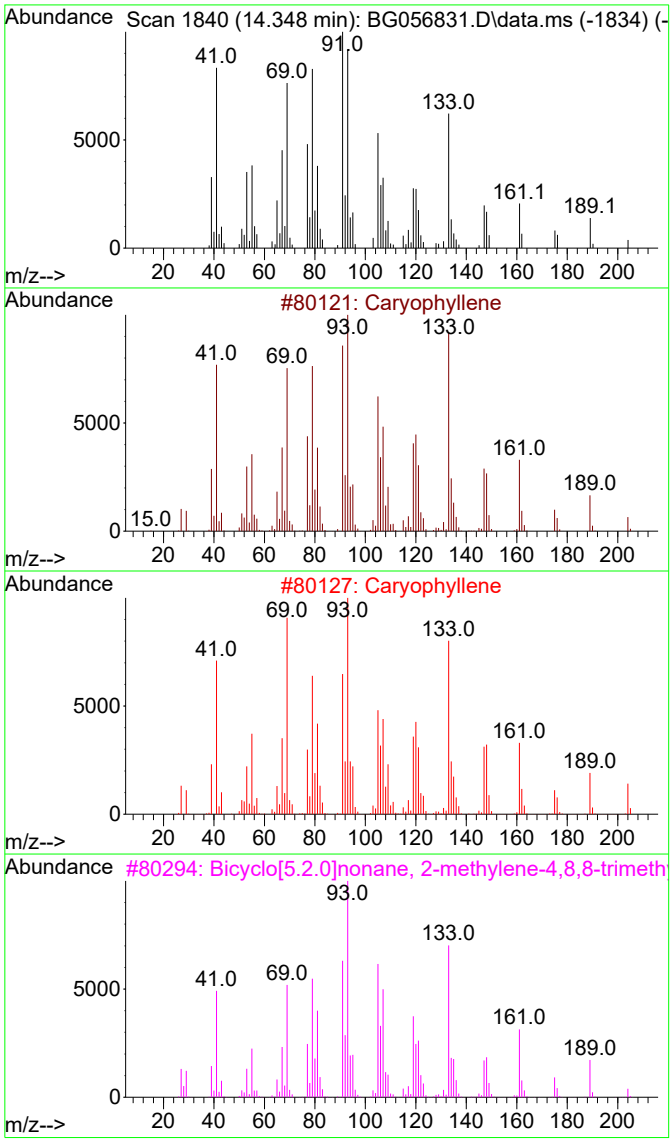
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

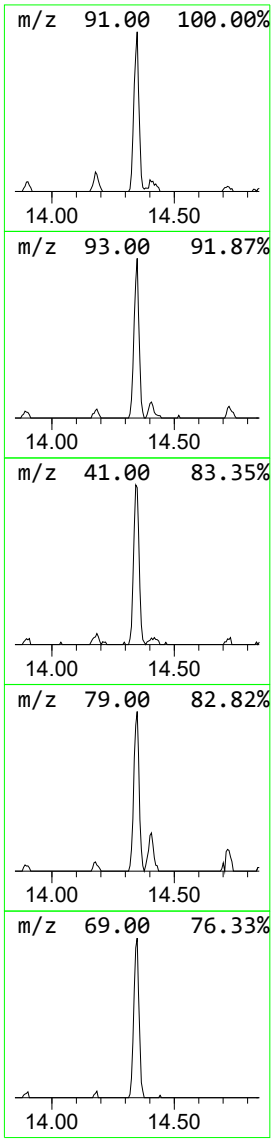
****

Peak Number 4 Caryophyllene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.348	22.53 ng/ul	259367	Acenaphthene-d10	15.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Caryophyllene	204	C15H24	000087-44-5	99
2	Caryophyllene	204	C15H24	000087-44-5	94
3	Bicyclo[5.2.0]nonane, 2-methylen...	204	C15H24	242794-76-9	91
4	Caryophyllene	204	C15H24	000087-44-5	87
5	Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	83





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Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

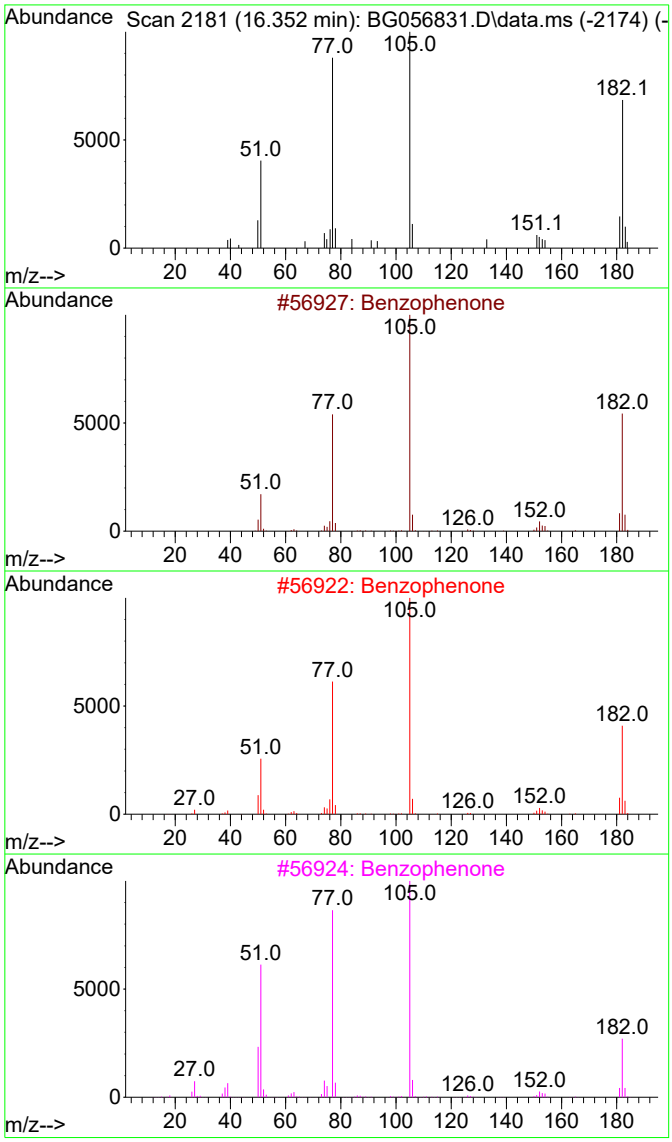
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

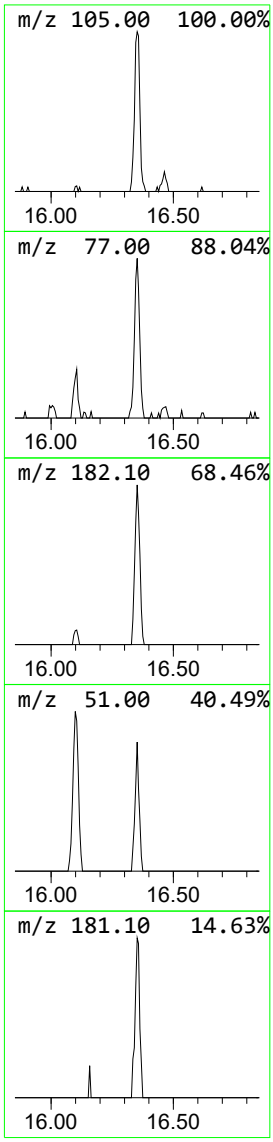
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.352	3.07 ng/ul	35403	Acenaphthene-d10	15.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	95
2	Benzophenone	182	C13H10O	000119-61-9	93
3	Benzophenone	182	C13H10O	000119-61-9	91
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	90





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ClientSampleId :
DBAD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

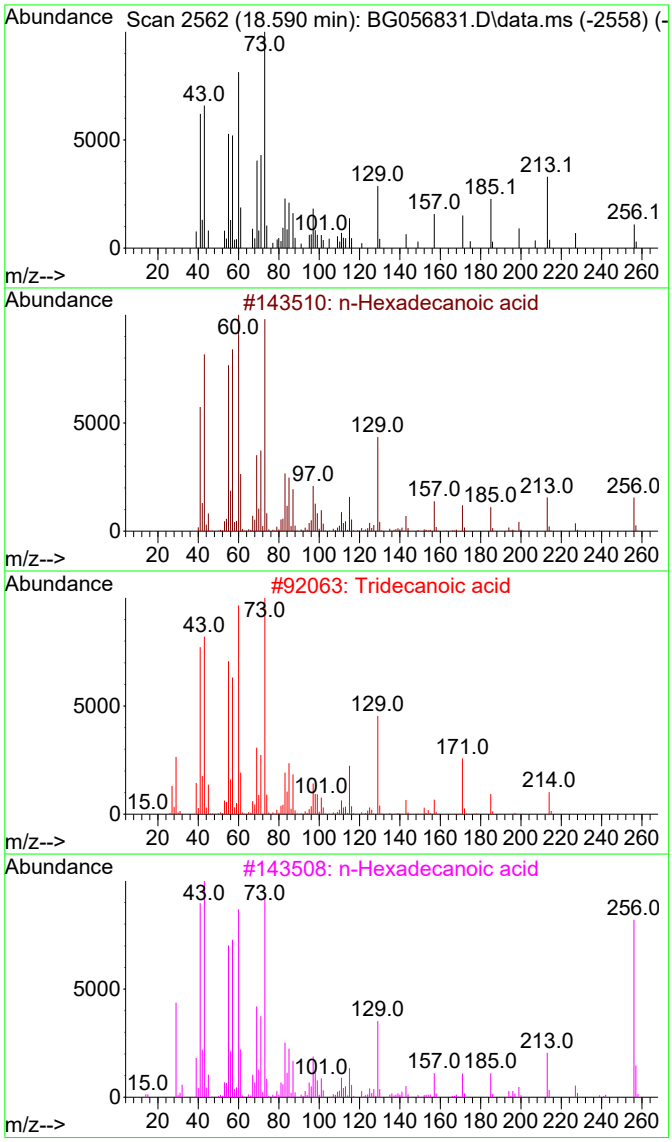
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

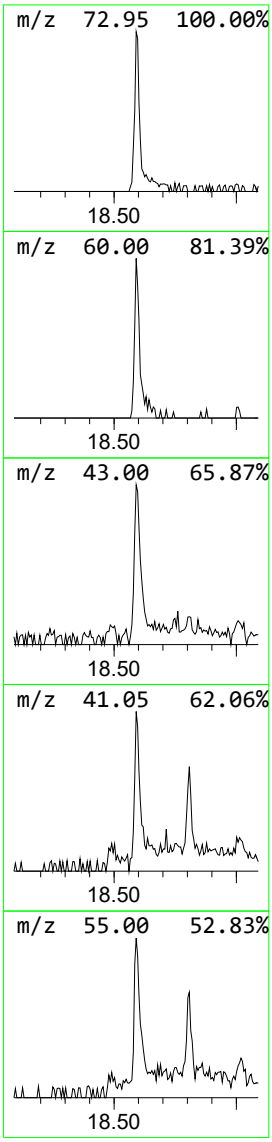
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.590	5.78 ng/ul	89859	Phenanthrene-d10	17.750

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	91
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056831.D
Acq On : 4 Mar 2023 10:24
Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

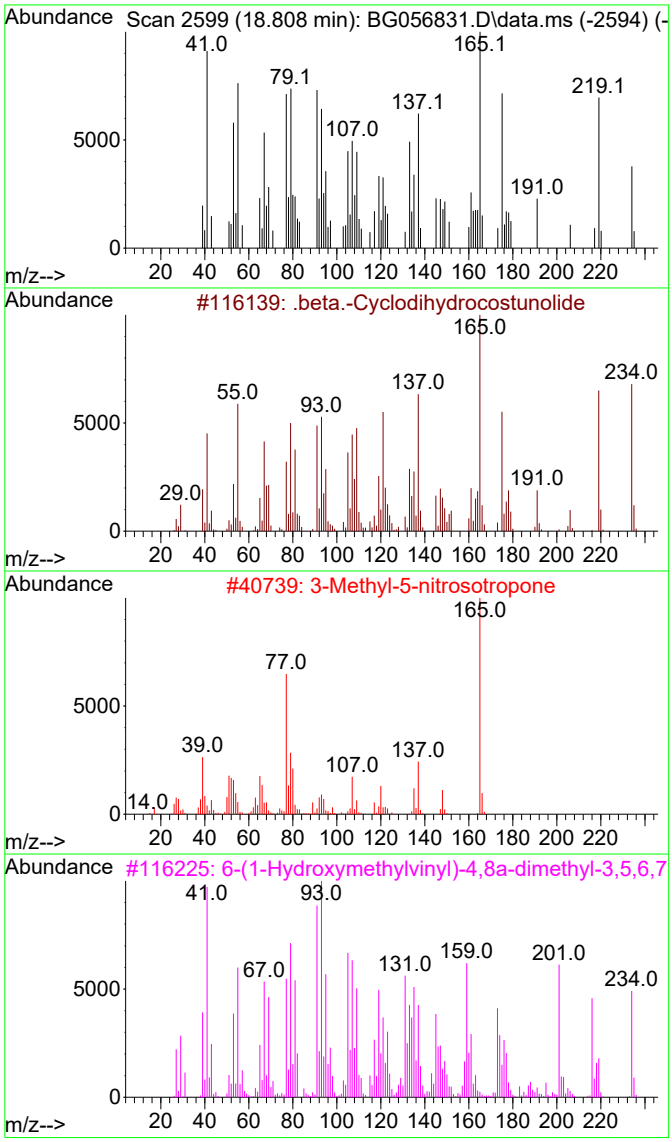
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

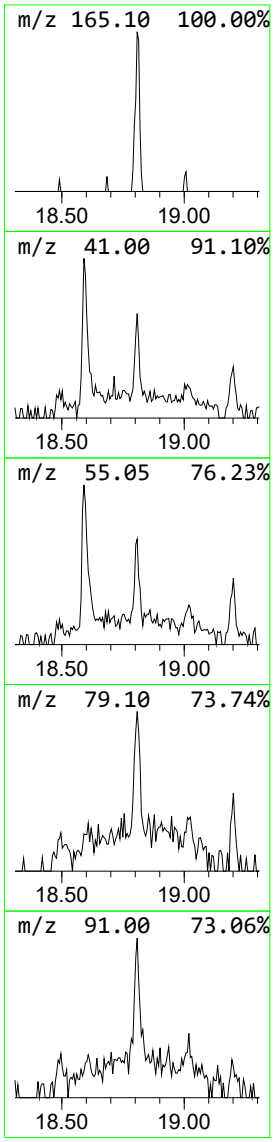
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 .beta.-Cyclodihydrocostunolide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.808	3.72 ng/ul	57834	Phenanthrene-d10	17.750

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Cyclodihydrocostunolide	234	C15H22O2	022387-64-0	95
2	3-Methyl-5-nitrosotropone	165	C8H7NO3	1000242-61-8	43
3	6-(1-Hydroxymethylvinyl)-4,8a-di...	234	C15H22O2	1000190-51-4	42
4	Quinoline, 2,3,4,4a,5,6-hexahydr...	165	C10H15NO	019500-64-2	38
5	1,3-Bis-(2-cyclopropyl,2-methylc...	258	C18H26O	1000222-08-6	35





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Data File : BG056831.D
Acq On : 4 Mar 2023 10:24
Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

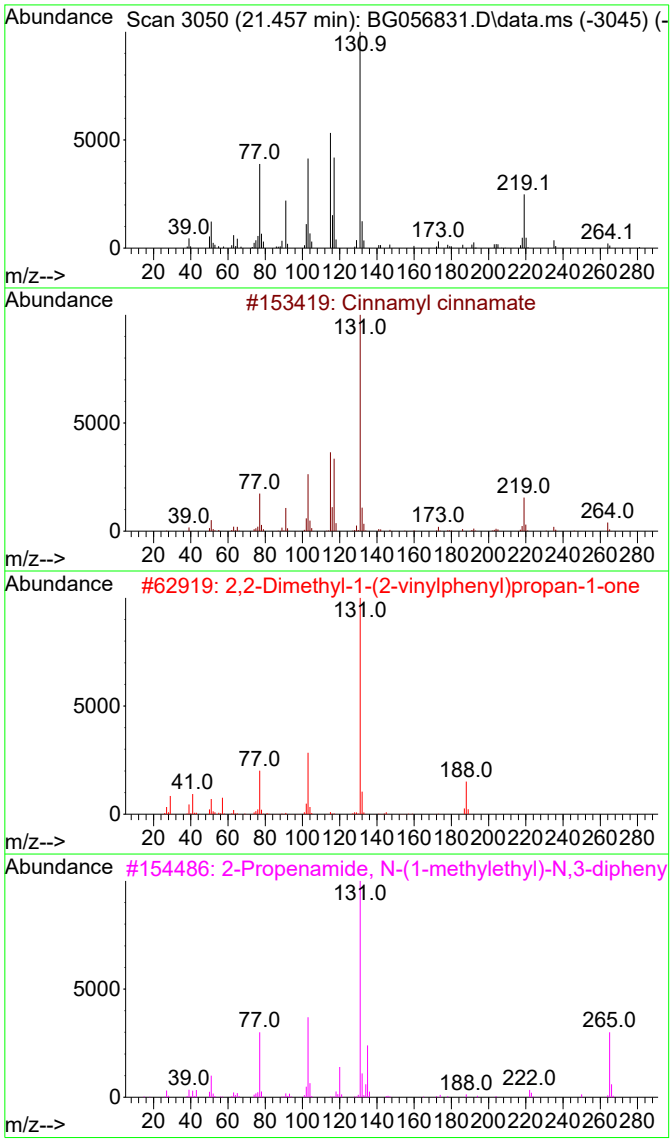
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

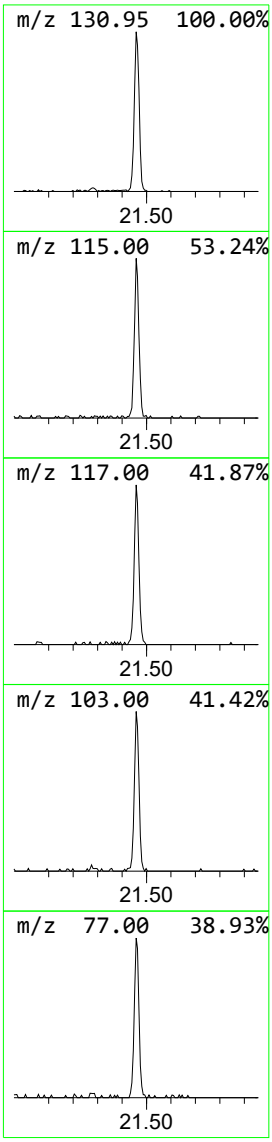
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cinnamyl cinnamate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.457	12.26 ng/ul	208273	Chrysene-d12	22.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	46
3	2-Propenamide, N-(1-methylethyl)...	265	C18H19NO	055044-37-6	46
4	p-Vinylbenzohydrazide	162	C9H10N2O	1000244-74-9	43
5	4-Methylcoumarine-7-cinnamate	306	C19H14O4	300829-05-4	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056831.D
Acq On : 4 Mar 2023 10:24
Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

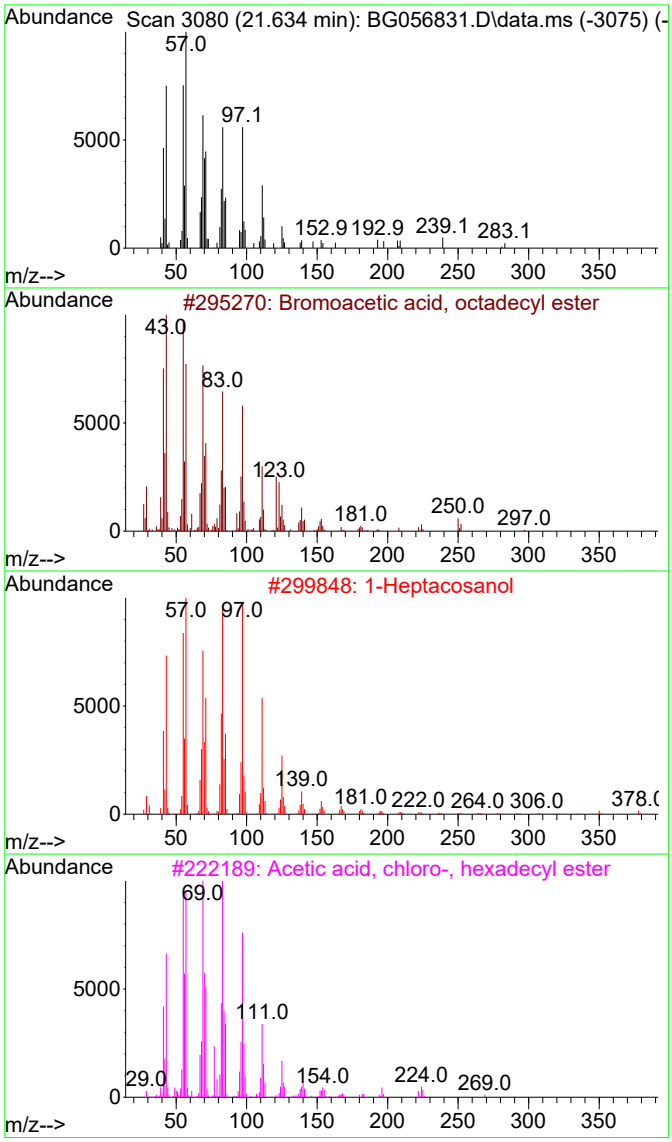
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

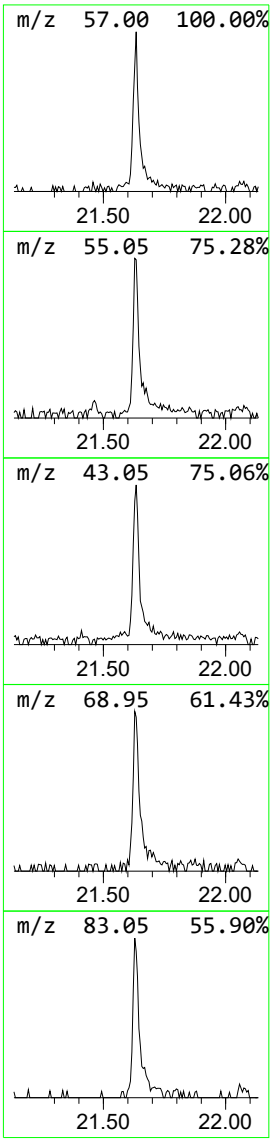
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Bromoacetic acid, octadecyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.634	6.42 ng/ul	109003	Chrysene-d12	22.057

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
2	1-Heptacosanol	396	C27H56O	002004-39-9	90
3	Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	87
4	2- Bromopropionic acid, pentadec...	362	C18H35BrO2	1000292-44-2	86
5	1-Hexacosanol	382	C26H54O	000506-52-5	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030323\
Data File : BG056831.D
Acq On : 4 Mar 2023 10:24
Operator : CG/JU
Sample : 01711-20
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBAD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(1S)-2,6,6-Trim...	7.080	2.9	ng/ul	13493	1	8.420	92584	20.0
Eucalyptol	8.725	5.0	ng/ul	22950	1	8.420	92584	20.0
1H-Cyclopropa[a...	14.178	2.1	ng/ul	24392	3	15.018	230275	20.0
Caryophyllene	14.348	22.5	ng/ul	259367	3	15.018	230275	20.0
Benzophenone	16.352	3.1	ng/ul	35403	3	15.018	230275	20.0
n-Hexadecanoic ...	18.590	5.8	ng/ul	89859	4	17.750	311137	20.0
.beta.-Cyclodih...	18.808	3.7	ng/ul	57834	4	17.750	311137	20.0
Cinnamyl cinnamate	21.457	12.3	ng/ul	208273	5	22.057	339626	20.0
Bromoacetic aci...	21.634	6.4	ng/ul	109003	5	22.057	339626	20.0