

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031822\
 Data File : BG052711.D
 Acq On : 18 Mar 2022 15:44
 Operator : CG/JU
 Sample : N1882-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBRY0

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-BG031822.M
 Title : SVOA CALIBRATION

Signal : TIC: BG052711.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.977	96	100	110	rBV	99549	181787	18.84%	1.815%
2	4.318	155	158	162	rBV4	4278	7158	0.74%	0.071%
3	5.235	312	314	318	rBV4	2266	3058	0.32%	0.031%
4	5.370	335	337	340	rBV3	1554	1827	0.19%	0.018%
5	5.722	395	397	400	rVB2	2500	2308	0.24%	0.023%
6	6.127	463	466	467	rBV3	1294	1502	0.16%	0.015%
7	6.380	507	509	513	rBV4	1522	2256	0.23%	0.023%
8	6.480	521	526	528	rVB2	1561	2073	0.21%	0.021%
9	6.562	537	540	542	rVB3	1797	2199	0.23%	0.022%
10	6.874	591	593	597	rVB4	1754	1768	0.18%	0.018%
11	6.938	603	604	609	rBV4	1207	1642	0.17%	0.016%
12	7.150	636	640	642	rVB3	1168	1688	0.17%	0.017%
13	7.291	657	664	674	rBV	153201	255160	26.45%	2.548%
14	7.402	679	683	686	rBV4	1122	1571	0.16%	0.016%
15	7.467	686	694	703	rBV	101260	175781	18.22%	1.755%
16	7.678	721	730	738	rVB	132027	230337	23.87%	2.300%
17	7.872	761	763	767	rVB3	1044	1395	0.14%	0.014%
18	8.148	803	810	818	rBV	132557	229516	23.79%	2.292%
19	8.207	818	820	827	rVB3	1567	2034	0.21%	0.020%
20	8.636	890	893	896	rBV3	1029	1378	0.14%	0.014%
21	8.783	916	918	921	rBV	968	1298	0.13%	0.013%
22	8.841	921	928	936	rVB	168724	295570	30.64%	2.952%
23	9.306	1000	1007	1014	rBV	126591	227161	23.54%	2.268%
24	9.729	1077	1079	1080	rBV2	1346	1297	0.13%	0.013%
25	9.758	1080	1084	1087	rVB3	5121	7075	0.73%	0.071%
26	9.864	1097	1102	1103	rBV3	974	1508	0.16%	0.015%
27	9.958	1117	1118	1124	rVB3	1037	1772	0.18%	0.018%
28	10.034	1124	1131	1137	rBV3	93632	169668	17.59%	1.694%
29	10.504	1207	1211	1215	rBV3	1329	1374	0.14%	0.014%
30	10.574	1215	1223	1230	rBV	162748	296463	30.73%	2.960%
31	10.733	1244	1250	1252	rBV3	2183	3280	0.34%	0.033%
32	10.833	1266	1267	1271	rVB	1664	1341	0.14%	0.013%
33	10.962	1282	1289	1297	rVB	174339	320095	33.18%	3.196%
34	11.085	1300	1310	1318	rBV	158501	295910	30.67%	2.955%
35	11.937	1453	1455	1459	rVB2	1026	1307	0.14%	0.013%
36	12.049	1470	1474	1477	rBV2	1187	1534	0.16%	0.015%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
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37	12.372	1522	1529	1533	rBV4	4832	6938	0.72%	0.069%
38	12.536	1551	1557	1562	rVB4	2969	4718	0.49%	0.047%
39	13.318	1687	1690	1692	rBV3	1432	1533	0.16%	0.015%
40	13.600	1734	1738	1743	rBV3	8466	12273	1.27%	0.123%
41	13.899	1782	1789	1791	rBV3	1101	2067	0.21%	0.021%
42	13.923	1791	1793	1798	rVV	1374	1779	0.18%	0.018%
43	14.164	1828	1834	1841	rVB	311852	455076	47.17%	4.544%
44	14.258	1849	1850	1853	rBV2	1274	1373	0.14%	0.014%
45	14.405	1872	1875	1878	rBV4	1617	2407	0.25%	0.024%
46	14.463	1878	1885	1893	rVB	326906	537855	55.75%	5.371%
47	14.663	1915	1919	1922	rBV3	2503	3389	0.35%	0.034%
48	14.769	1930	1937	1945	rVB	283556	453234	46.98%	4.526%
49	14.927	1958	1964	1972	rBV	131454	212285	22.00%	2.120%
50	15.568	2071	2073	2078	rVB5	1918	2422	0.25%	0.024%
51	15.621	2078	2082	2084	rBV2	1117	1510	0.16%	0.015%
52	15.756	2098	2105	2113	rVB	434971	688651	71.38%	6.877%
53	15.856	2117	2122	2131	rVV	99926	147756	15.31%	1.475%
54	15.996	2142	2146	2149	rVV3	1920	2405	0.25%	0.024%
55	16.026	2149	2151	2158	rVB3	1048	1593	0.17%	0.016%
56	16.478	2225	2228	2229	rBV3	2136	1543	0.16%	0.015%
57	16.748	2273	2274	2279	rBV2	1872	2572	0.27%	0.026%
58	16.801	2281	2283	2287	rBV2	1308	1305	0.14%	0.013%
59	16.913	2300	2302	2304	rBV3	1818	1296	0.13%	0.013%
60	17.136	2337	2340	2343	rVB4	1698	1867	0.19%	0.019%
61	17.412	2385	2387	2391	rVB3	2391	2175	0.23%	0.022%
62	17.442	2391	2392	2395	rVB2	1835	1399	0.15%	0.014%
63	17.512	2397	2404	2409	rBV	383983	584625	60.60%	5.838%
64	17.612	2414	2421	2428	rVV	481548	772203	80.04%	7.711%
65	17.682	2430	2433	2438	rVV4	5205	6431	0.67%	0.064%
66	17.888	2466	2468	2471	rVB2	1743	1465	0.15%	0.015%
67	18.129	2506	2509	2511	rBV4	2106	2272	0.24%	0.023%
68	18.176	2515	2517	2520	rVB3	1736	1612	0.17%	0.016%
69	18.387	2549	2553	2557	rBV3	31114	40485	4.20%	0.404%
70	18.558	2579	2582	2583	rBV3	2067	2206	0.23%	0.022%
71	18.593	2586	2588	2591	rBV3	2048	2144	0.22%	0.021%
72	18.992	2654	2656	2657	rBV	3248	2162	0.22%	0.022%
73	19.462	2732	2736	2739	rVB4	6432	8077	0.84%	0.081%
74	19.521	2742	2746	2751	rBV3	10133	20011	2.07%	0.200%
75	19.656	2766	2769	2771	rBV4	7393	7912	0.82%	0.079%
76	19.891	2803	2809	2818	rBV	659727	964806	100.00%	9.635%

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Integrator: RTE
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Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Title : SVOA CALIBRATION

77	21.430	3067	3071	3079	rVB	78006	104967	10.88%	1.048%
78	21.795	3127	3133	3141	rVB	413814	691631	71.69%	6.907%
79	24.884	3651	3659	3671	rVB	313842	899212	93.20%	8.980%
80	25.119	3691	3699	3709	rVB2	218065	617269	63.98%	6.164%

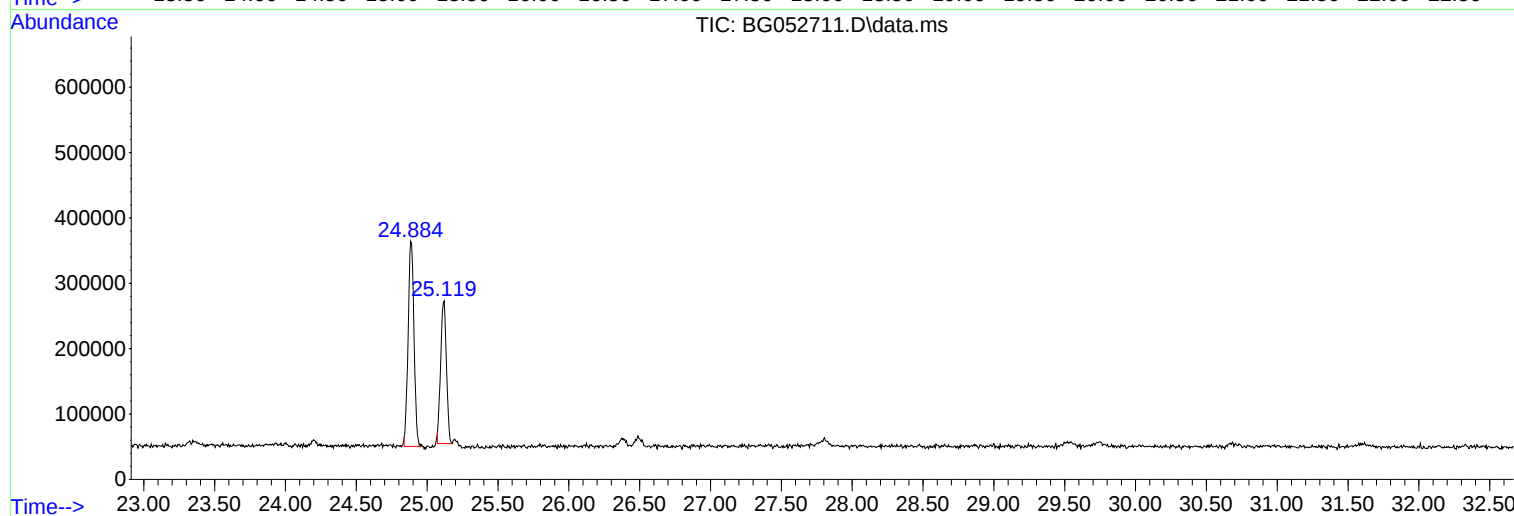
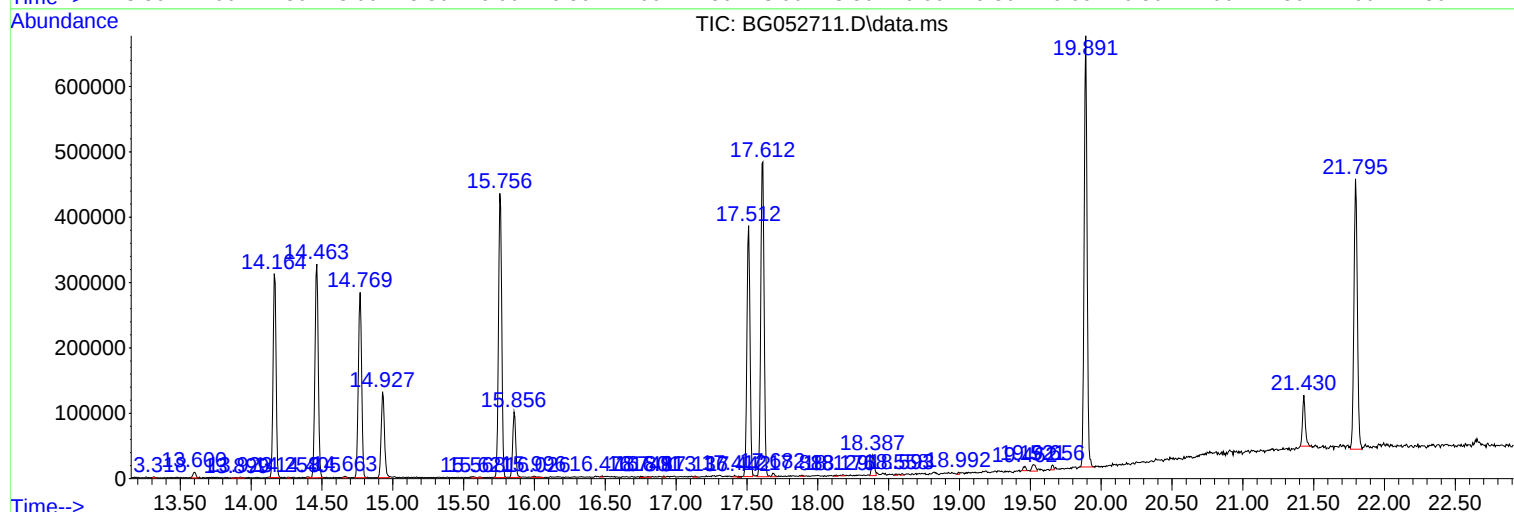
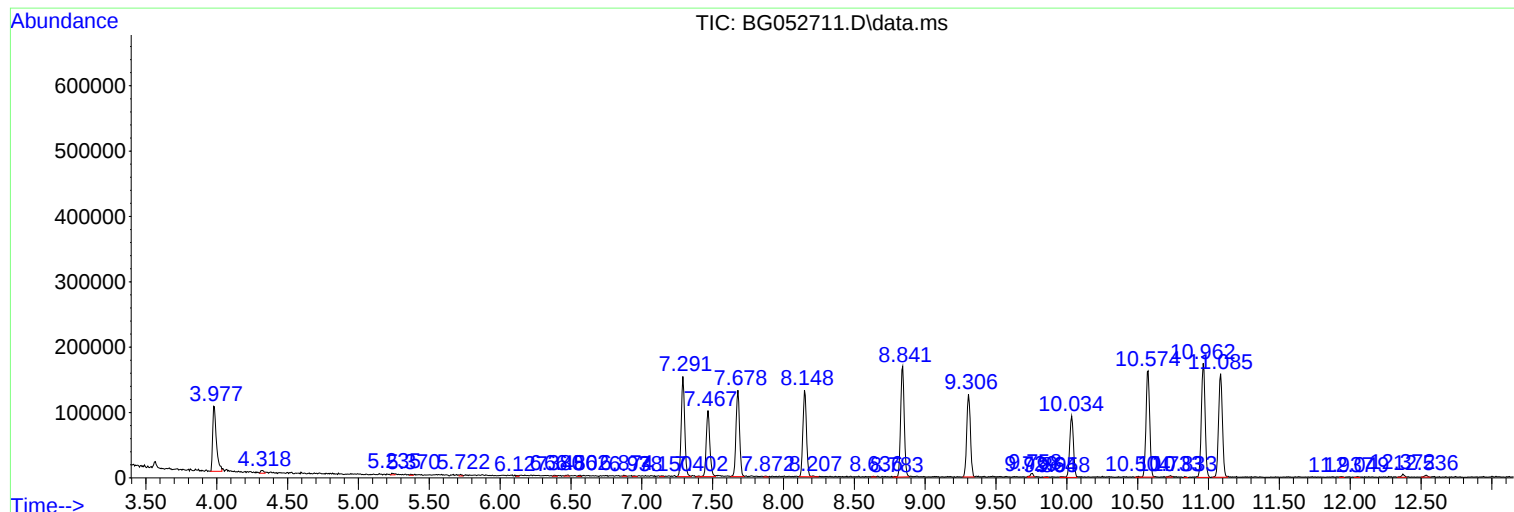
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION

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



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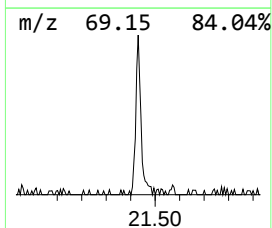
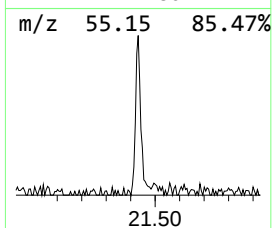
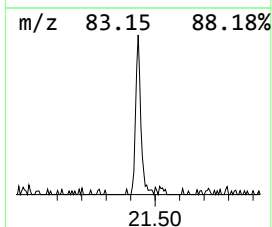
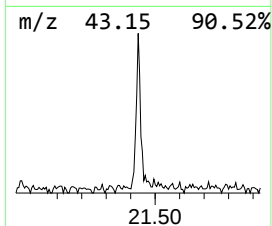
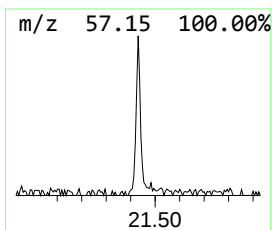
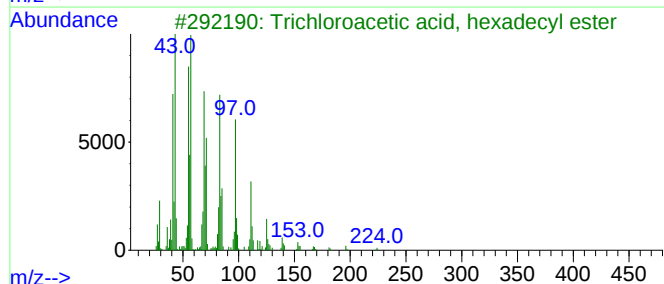
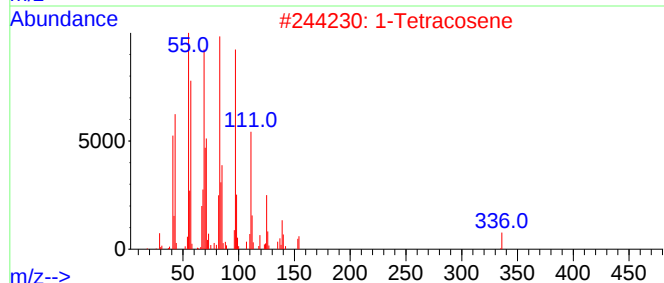
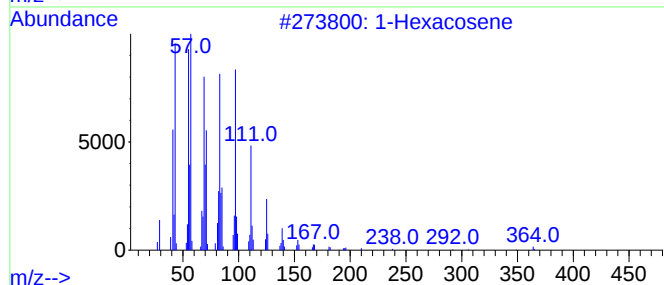
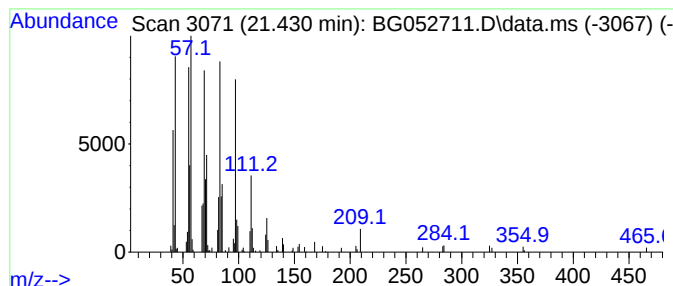
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.430	3.04 ng/ul	104967	Chrysene-d12	21.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	91
2	1-Tetracosene			336	C24H48	010192-32-2	90
3	Trichloroacetic acid, hexadecyl ...			386	C18H33Cl3O2	074339-54-1	87
4	Trifluoroacetoxy hexadecane			338	C18H33F3O2	006222-03-3	87
5	Pentacos-1-ene			350	C25H50	016980-85-1	87



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	21.430	3.0	ng/ul	104967	5	21.794	691631	20.0