

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122721\
 Data File : BG051815.D
 Acq On : 28 Dec 2021 00:06
 Operator : CG/JU
 Sample : M5193-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0J04

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-BG121421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG051815.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.601	15	19	24	rVB2	8607	13587	1.09%	0.123%
2	4.030	88	92	108	rBV	32019	63029	5.04%	0.572%
3	4.382	148	152	157	rVB2	2337	3812	0.30%	0.035%
4	4.935	242	246	251	rBV3	2821	4226	0.34%	0.038%
5	7.085	608	612	619	rBV2	997	2030	0.16%	0.018%
6	7.408	661	667	676	rBV	35660	66494	5.32%	0.604%
7	7.578	689	696	704	rBV	134634	230547	18.44%	2.094%
8	7.795	726	733	742	rBV	132778	227072	18.16%	2.062%
9	8.271	806	814	821	rVB	121750	213041	17.04%	1.935%
10	8.970	926	933	945	rVB	103115	184523	14.76%	1.676%
11	9.440	1004	1013	1023	rBV	192099	346907	27.75%	3.151%
12	9.863	1081	1085	1089	rVB2	3604	5267	0.42%	0.048%
13	10.175	1131	1138	1146	rBV	151921	270224	21.62%	2.454%
14	10.721	1223	1231	1242	rBV	210749	378809	30.30%	3.441%
15	11.109	1288	1297	1305	rBV	175839	313634	25.09%	2.849%
16	11.232	1310	1318	1328	rVV	175727	321344	25.70%	2.919%
17	12.366	1507	1511	1515	rVB3	2769	3365	0.27%	0.031%
18	12.495	1530	1533	1539	rBV4	1476	2717	0.22%	0.025%
19	12.683	1561	1565	1569	rVB3	3542	5392	0.43%	0.049%
20	13.118	1635	1639	1643	rBV2	979	1757	0.14%	0.016%
21	13.717	1739	1741	1748	rVB5	4454	6821	0.55%	0.062%
22	14.293	1833	1839	1846	rBV	437686	659411	52.75%	5.989%
23	14.527	1876	1879	1881	rBV2	2438	2593	0.21%	0.024%
24	14.604	1885	1892	1899	rVB	491094	776086	62.08%	7.049%
25	14.909	1937	1944	1952	rVB	277464	438831	35.10%	3.986%
26	15.080	1966	1973	1981	rBV2	20713	41062	3.28%	0.373%
27	15.526	2046	2049	2053	rBV	2167	3098	0.25%	0.028%
28	15.691	2072	2077	2082	rBV3	3196	6043	0.48%	0.055%
29	15.726	2082	2083	2088	rVB3	2921	3210	0.26%	0.029%
30	15.896	2104	2112	2122	rBV	641100	980464	78.43%	8.905%
31	15.996	2123	2129	2135	rVV	171181	259163	20.73%	2.354%
32	16.143	2151	2154	2157	rVB2	3212	3712	0.30%	0.034%
33	16.237	2167	2170	2173	rVB4	2200	2737	0.22%	0.025%
34	16.466	2205	2209	2210	rBV2	1615	1655	0.13%	0.015%
35	16.531	2217	2220	2222	rBV3	2406	2827	0.23%	0.026%
36	16.566	2224	2226	2228	rBV3	2938	2397	0.19%	0.022%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
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37	16.777	2258	2262	2265	rVB4	2696	3126	0.25%	0.028%
38	16.936	2285	2289	2292	rBV2	1308	2138	0.17%	0.019%
39	17.001	2297	2300	2303	rVB3	2097	2674	0.21%	0.024%
40	17.036	2303	2306	2308	rBV2	2885	3229	0.26%	0.029%
41	17.112	2313	2319	2322	rBV4	3108	4526	0.36%	0.041%
42	17.382	2361	2365	2371	rVB4	4609	6872	0.55%	0.062%
43	17.441	2371	2375	2379	rVB4	3080	4293	0.34%	0.039%
44	17.541	2388	2392	2398	rVB4	8125	8765	0.70%	0.080%
45	17.653	2404	2411	2419	rBV	364908	531055	42.48%	4.823%
46	17.753	2419	2428	2438	rVB	709308	1098217	87.85%	9.975%
47	17.917	2452	2456	2460	rBV3	4371	5219	0.42%	0.047%
48	18.029	2473	2475	2480	rVB4	1583	2022	0.16%	0.018%
49	18.416	2536	2541	2542	rBV4	1944	2585	0.21%	0.023%
50	18.522	2555	2559	2566	rBV5	7124	10502	0.84%	0.095%
51	18.593	2566	2571	2577	rVV3	8439	13804	1.10%	0.125%
52	18.663	2581	2583	2586	rVB2	2471	2133	0.17%	0.019%
53	18.845	2611	2614	2615	rVB3	2460	1978	0.16%	0.018%
54	18.939	2626	2630	2634	rVB5	5782	8903	0.71%	0.081%
55	19.133	2657	2663	2666	rBV5	7337	10462	0.84%	0.095%
56	19.221	2676	2678	2683	rBV4	1769	2577	0.21%	0.023%
57	19.262	2683	2685	2687	rVV3	2105	1917	0.15%	0.017%
58	19.386	2703	2706	2708	rVB3	5665	6379	0.51%	0.058%
59	19.597	2738	2742	2745	rBV2	9759	13551	1.08%	0.123%
60	19.668	2749	2754	2756	rBV	12093	16381	1.31%	0.149%
61	19.697	2757	2759	2762	rVB4	6783	5173	0.41%	0.047%
62	19.903	2791	2794	2795	rBV3	2705	2804	0.22%	0.025%
63	20.032	2809	2816	2822	rVB	856704	1250166	100.00%	11.355%
64	21.577	3075	3079	3085	rBV4	39149	68967	5.52%	0.626%
65	21.976	3140	3147	3155	rVB	307140	538765	43.10%	4.893%
66	25.231	3690	3701	3714	rBV2	331696	1029317	82.33%	9.349%
67	25.472	3732	3742	3759	rVB2	145092	503773	40.30%	4.576%

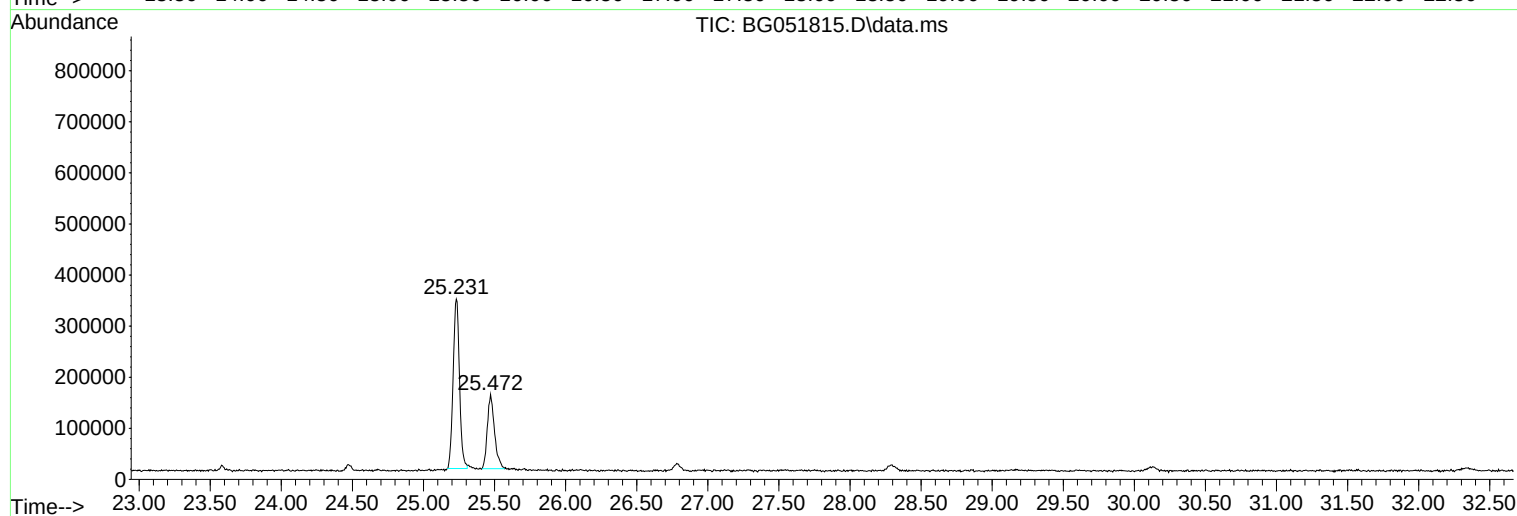
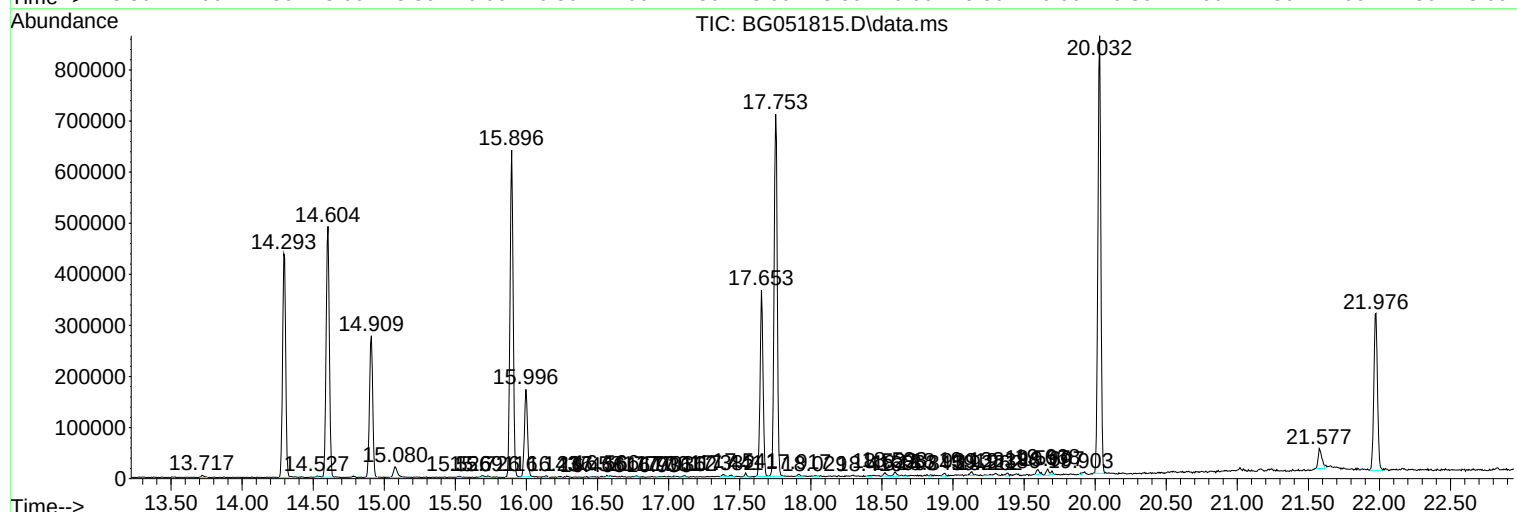
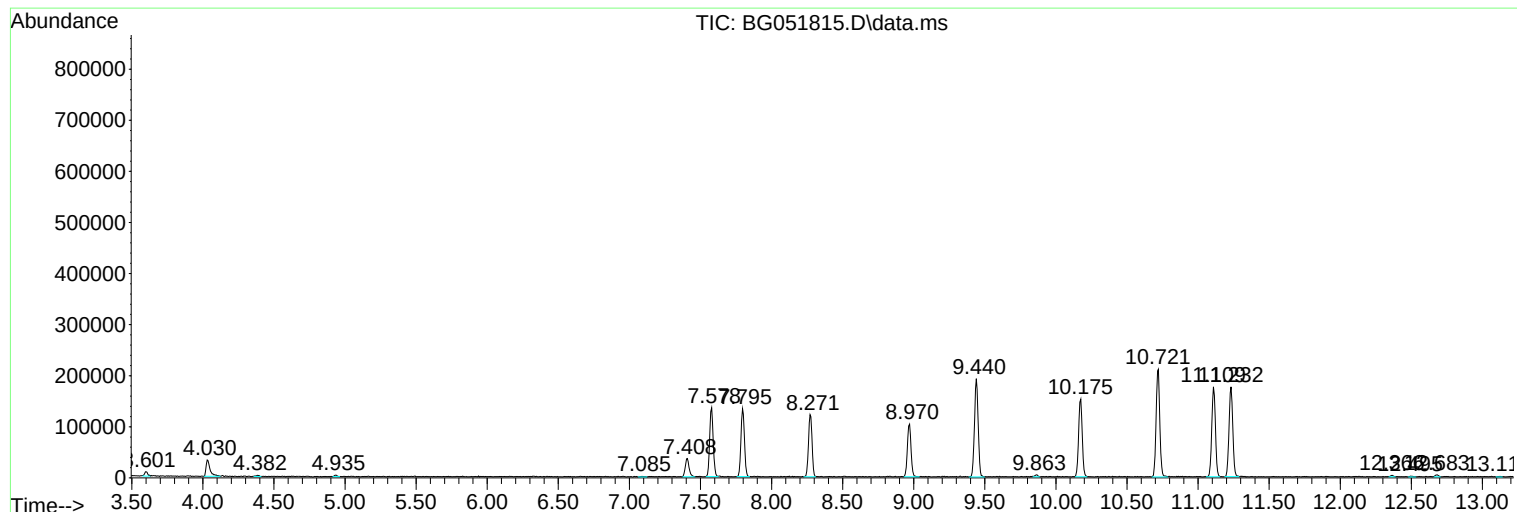
Sum of corrected areas: 11010160

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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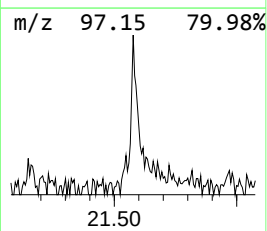
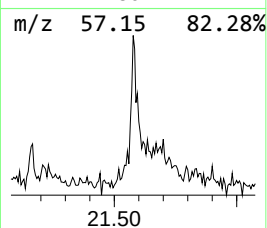
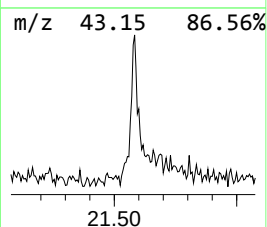
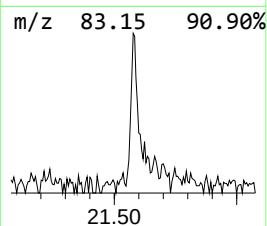
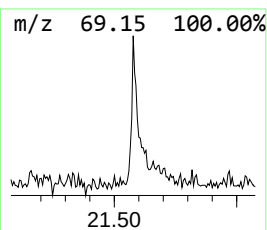
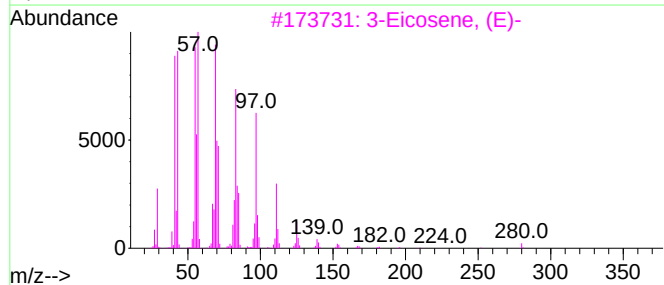
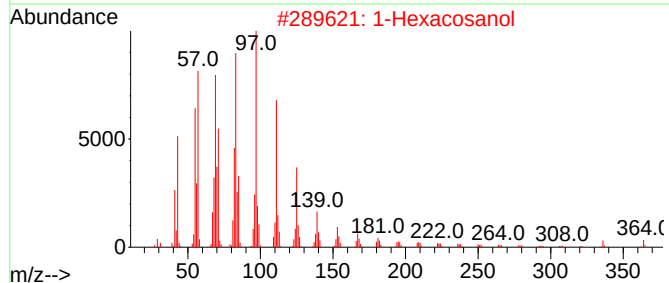
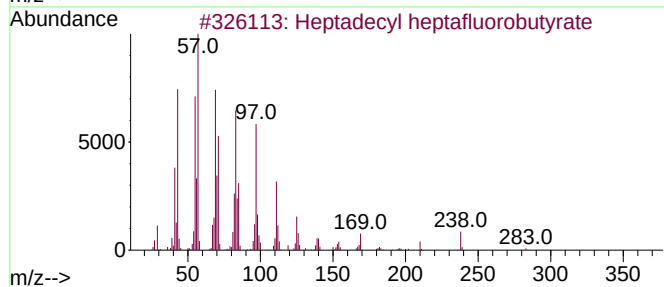
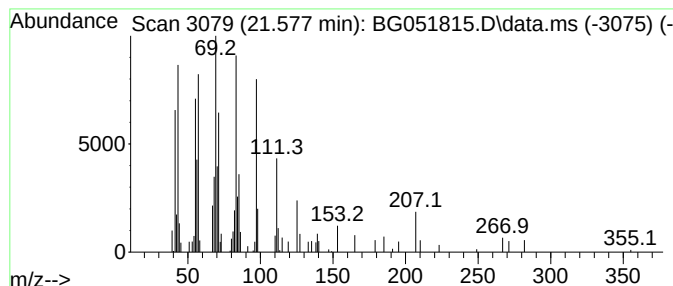
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Heptadecyl heptafluorobutyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.577	2.56 ng/ul	68967	Chrysene-d12	21.970

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
2			1-Hexacosanol	382	C26H54O	000506-52-5	91
3			3-Eicosene, (E)-	280	C20H40	074685-33-9	87
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	83
5			Behenic alcohol	326	C22H46O	000661-19-8	81



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Heptadecyl hept...	21.577	2.6	ng/ul	68967	5	21.970	538765	20.0