

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050823\
 Data File : BP014998.D
 Acq On : 08 May 2023 17:52
 Operator : CG/JU
 Sample : 02589-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JRECO

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_P\Methods\SFAM-EPA-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014998.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.458	9	12	19	rVB	33320	48735	0.98%	0.099%
2	3.946	94	95	97	rBV2	9688	9427	0.19%	0.019%
3	4.146	124	129	134	rVB2	18762	31582	0.63%	0.064%
4	4.246	142	146	150	rVB	17713	26539	0.53%	0.054%
5	5.199	303	308	315	rVB	14988	25933	0.52%	0.052%
6	5.693	390	392	398	rVB6	5824	6938	0.14%	0.014%
7	5.817	409	413	419	rVB	8288	12090	0.24%	0.024%
8	7.305	659	666	680	rBV	525635	879504	17.65%	1.778%
9	7.481	689	696	705	rVB	491065	774727	15.55%	1.566%
10	7.687	723	731	740	rBV	571358	913804	18.34%	1.847%
11	7.769	741	745	750	rVB4	9994	13303	0.27%	0.027%
12	8.164	805	812	819	rBV	1197349	1881984	37.77%	3.805%
13	8.869	926	932	946	rBV	647357	1083239	21.74%	2.190%
14	8.999	948	954	962	rVB	32701	54785	1.10%	0.111%
15	9.340	1004	1012	1027	rBV	558801	928740	18.64%	1.878%
16	9.499	1034	1039	1041	rVB6	3930	5354	0.11%	0.011%
17	9.740	1076	1080	1085	rVB2	11274	15768	0.32%	0.032%
18	10.069	1128	1136	1151	rBV	444219	758368	15.22%	1.533%
19	10.428	1195	1197	1203	rVB7	4443	5686	0.11%	0.011%
20	10.610	1216	1228	1238	rBV	702036	1205263	24.19%	2.437%
21	10.999	1286	1294	1307	rBV	1806267	3032449	60.85%	6.130%
22	11.134	1310	1317	1329	rBV	522650	961256	19.29%	1.943%
23	11.528	1379	1384	1388	rBV7	4939	7977	0.16%	0.016%
24	12.399	1526	1532	1536	rBV5	7815	14310	0.29%	0.029%
25	12.481	1541	1546	1551	rVB4	5619	9936	0.20%	0.020%
26	12.575	1557	1562	1567	rBV2	11448	20195	0.41%	0.041%
27	13.334	1685	1691	1697	rVB10	4254	7637	0.15%	0.015%
28	13.616	1733	1739	1744	rBV	13767	21558	0.43%	0.044%
29	13.693	1746	1752	1757	rBV	43215	67218	1.35%	0.136%
30	13.781	1762	1767	1770	rVB7	4035	6721	0.13%	0.014%
31	14.116	1822	1824	1831	rVB7	3822	7724	0.16%	0.016%
32	14.204	1831	1839	1854	rBV	1322996	1891524	37.96%	3.824%
33	14.328	1854	1860	1864	rBV9	5325	9397	0.19%	0.019%
34	14.434	1875	1878	1880	rBV4	4282	5491	0.11%	0.011%
35	14.504	1882	1890	1900	rBV	1638625	2424466	48.65%	4.901%
36	14.687	1916	1921	1925	rVB7	5778	8294	0.17%	0.017%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Title : SVOA CALIBRATION

37	14.810	1934	1942	1952	rBV	2914076	4264198	85.57%	8.620%
38	14.993	1966	1973	1988	rBV	255049	478740	9.61%	0.968%
39	15.216	2006	2011	2015	rVB8	12807	21965	0.44%	0.044%
40	15.798	2101	2110	2123	rBV	2087344	2951264	59.22%	5.966%
41	15.916	2124	2130	2142	rVV	449622	653138	13.11%	1.320%
42	16.040	2146	2151	2161	rVB5	13151	27924	0.56%	0.056%
43	16.157	2166	2171	2178	rBV	184689	254953	5.12%	0.515%
44	16.498	2226	2229	2235	rBV8	4688	7887	0.16%	0.016%
45	16.587	2240	2244	2248	rVB6	5134	7026	0.14%	0.014%
46	16.945	2301	2305	2307	rBV5	5359	7642	0.15%	0.015%
47	17.298	2361	2365	2368	rBV6	4966	7221	0.14%	0.015%
48	17.351	2370	2374	2379	rVB4	9553	14403	0.29%	0.029%
49	17.434	2385	2388	2390	rBV4	6125	7278	0.15%	0.015%
50	17.510	2397	2401	2403	rBV5	5958	8267	0.17%	0.017%
51	17.563	2403	2410	2420	rVV	3507118	4983192	100.00%	10.074%
52	17.663	2420	2427	2437	rBV	2194384	3080830	61.82%	6.228%
53	18.310	2533	2537	2543	rVB8	7650	11849	0.24%	0.024%
54	18.416	2550	2555	2565	rBV	318613	483548	9.70%	0.978%
55	19.304	2703	2706	2710	rBV5	16123	21985	0.44%	0.044%
56	19.451	2729	2731	2736	rVB3	26108	35809	0.72%	0.072%
57	19.639	2759	2763	2772	rBV3	100122	173179	3.48%	0.350%
58	19.781	2785	2787	2793	rBV3	24296	30140	0.60%	0.061%
59	19.881	2798	2804	2813	rBV	2699495	3561699	71.47%	7.200%
60	20.375	2886	2888	2892	rVB	55291	62027	1.24%	0.125%
61	20.857	2967	2970	2974	rBV	97985	112096	2.25%	0.227%
62	21.316	3044	3048	3056	rBV	620368	829178	16.64%	1.676%
63	21.645	3099	3104	3113	rBV	3049313	4181780	83.92%	8.454%
64	21.775	3123	3126	3132	rVB2	126470	145541	2.92%	0.294%
65	24.074	3511	3517	3531	rVB2	1032732	2475190	49.67%	5.004%
66	24.245	3539	3546	3558	rVB2	1530707	3388567	68.00%	6.850%

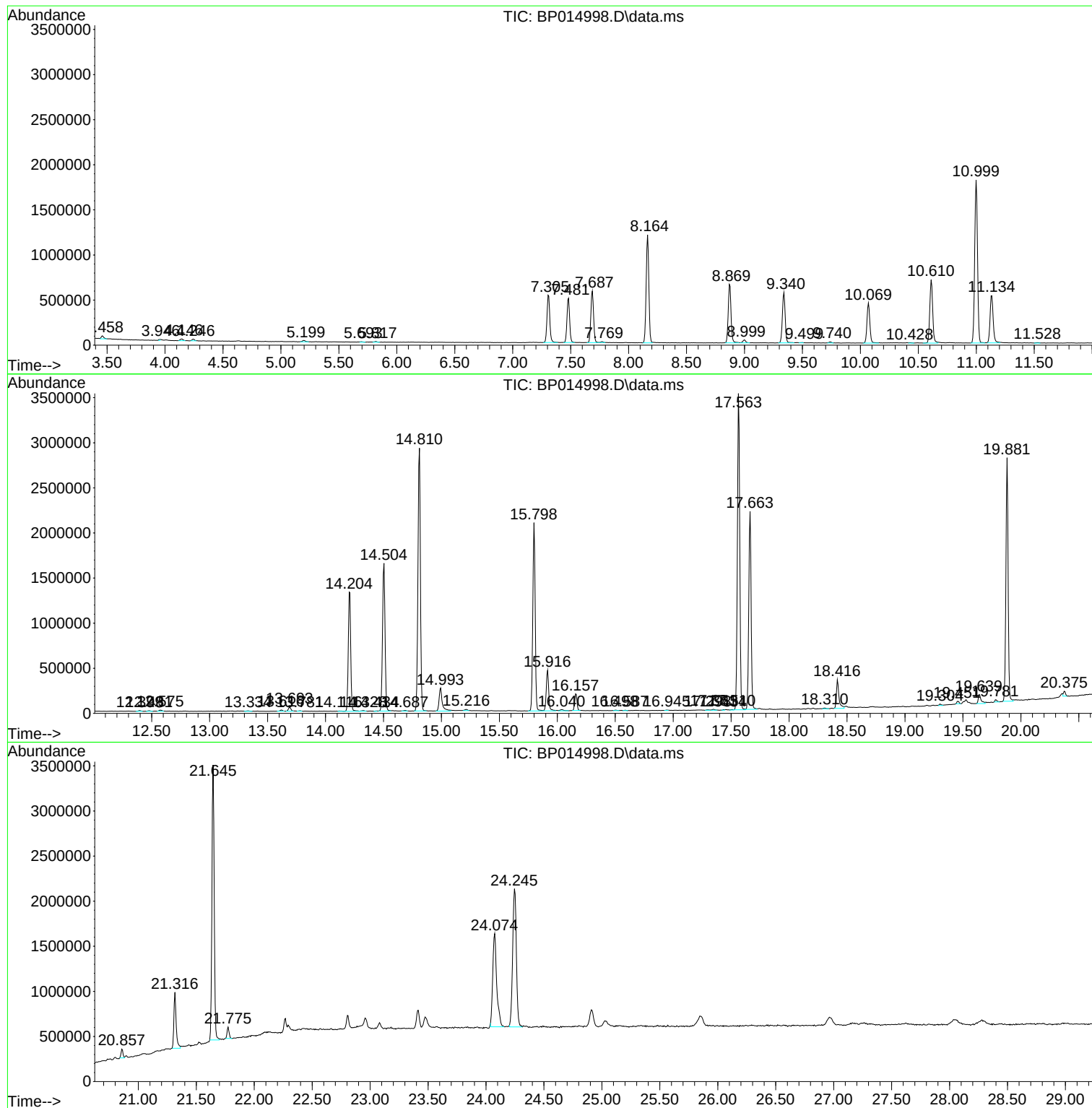
Sum of corrected areas: 49466438

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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



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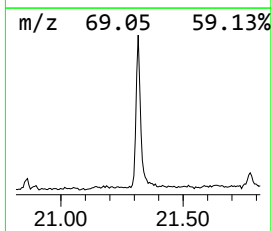
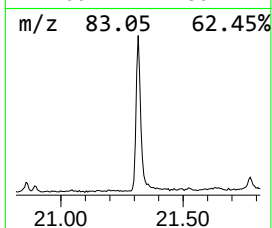
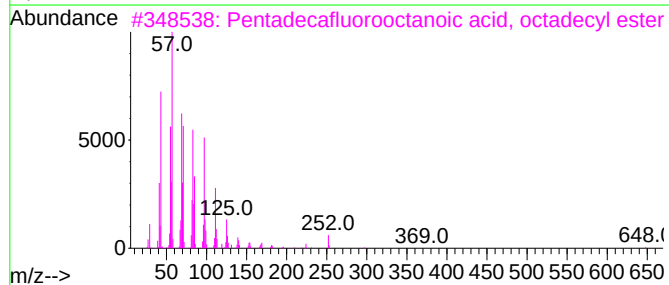
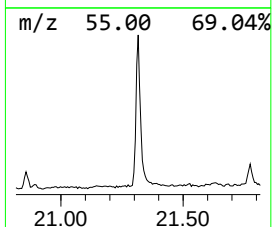
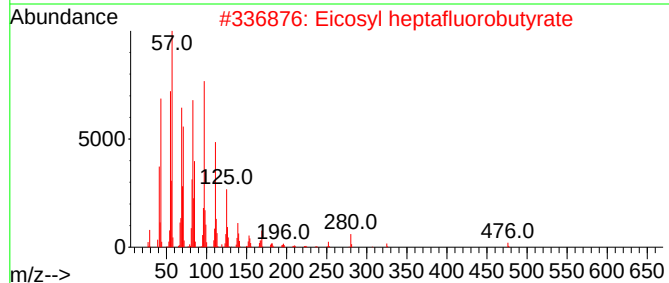
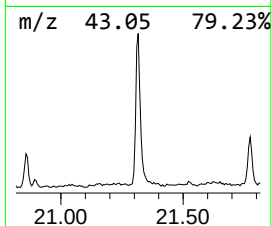
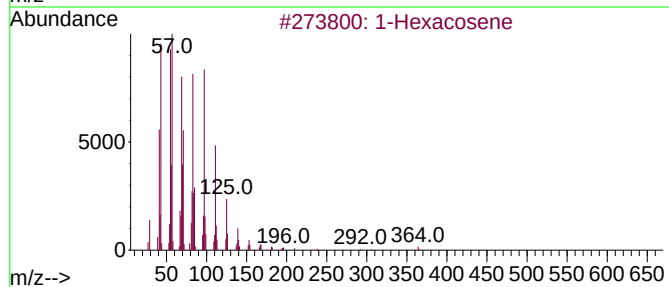
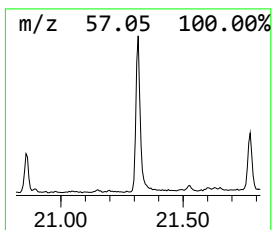
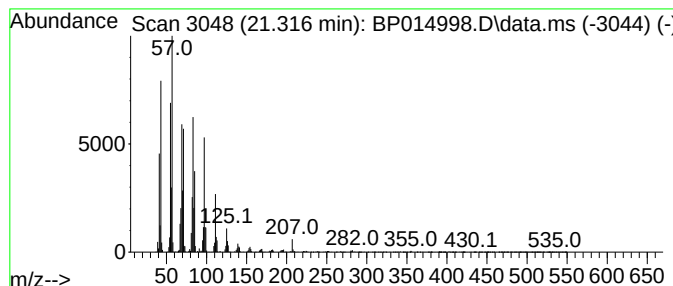
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

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.
21.316	3.97 ng/ul	829178	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	93
2	Eicosyl heptafluorobutyrate			494	C24H41F7O2	1000351-83-5	91
3	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	91
4	Nonadecyl pentafluoropropionate			430	C22H39F5O2	1000351-88-8	91
5	Cyclotetradecane			196	C14H28	000295-17-0	90



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	21.316	4.0	ng/u1	829178	5	21.645	4181780	20.0