

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051024\
 Data File : BP020301.D
 Acq On : 10 May 2024 22:16
 Operator : MA/JU
 Sample : P2370-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43P7

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020301.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.216	19	22	27	rVB3	13400	17121	0.97%	0.092%
2	3.622	88	91	111	rVB	87761	182458	10.35%	0.985%
3	3.905	137	139	149	rVB5	5441	9184	0.52%	0.050%
4	4.799	289	291	292	rBV2	2555	1820	0.10%	0.010%
5	6.710	612	616	619	rBV6	1273	2232	0.13%	0.012%
6	6.810	627	633	653	rBV	176498	406943	23.08%	2.197%
7	6.969	655	660	678	rVB	152890	287755	16.32%	1.553%
8	7.157	684	692	708	rBV	207977	409274	23.21%	2.209%
9	7.616	762	770	788	rBV	269570	470710	26.69%	2.541%
10	8.328	884	891	910	rBV	221726	464727	26.35%	2.508%
11	8.775	959	967	978	rBV	203306	412997	23.42%	2.229%
12	9.116	1021	1025	1028	rBV5	2394	3560	0.20%	0.019%
13	9.375	1064	1069	1072	rBV7	2927	5779	0.33%	0.031%
14	9.493	1082	1089	1103	rBV	183209	366755	20.80%	1.980%
15	10.028	1171	1180	1202	rBV	181508	538412	30.53%	2.906%
16	10.251	1212	1218	1227	rBV8	6034	16444	0.93%	0.089%
17	10.387	1233	1241	1252	rBV	356517	653774	37.07%	3.529%
18	10.551	1262	1269	1296	rBV	205133	525830	29.82%	2.838%
19	10.851	1317	1320	1322	rBV4	1785	2118	0.12%	0.011%
20	11.198	1373	1379	1381	rBV6	6338	11896	0.67%	0.064%
21	12.010	1509	1517	1523	rBV8	3687	8802	0.50%	0.048%
22	13.087	1694	1700	1705	rVB9	1367	3102	0.18%	0.017%
23	13.163	1707	1713	1724	rVV3	9572	17740	1.01%	0.096%
24	13.251	1724	1728	1732	rVB7	1503	2407	0.14%	0.013%
25	13.704	1797	1805	1824	rBV	470048	879925	49.90%	4.749%
26	13.881	1833	1835	1839	rVB5	1515	1800	0.10%	0.010%
27	13.975	1843	1851	1861	rBV	650968	1017966	57.72%	5.495%
28	14.186	1884	1887	1891	rVB5	1536	1789	0.10%	0.010%
29	14.286	1896	1904	1918	rBV	541908	917697	52.04%	4.953%
30	14.522	1932	1944	1948	rBV7	20270	47651	2.70%	0.257%
31	14.586	1948	1955	1965	rBV3	75182	235699	13.37%	1.272%
32	15.045	2030	2033	2035	rBV4	1595	1867	0.11%	0.010%
33	15.110	2042	2044	2047	rVB4	2264	2369	0.13%	0.013%
34	15.310	2070	2078	2094	rBV	734989	1280381	72.60%	6.911%
35	15.475	2099	2106	2117	rBV	146231	259482	14.71%	1.401%
36	15.892	2174	2177	2178	rBV3	2145	2216	0.13%	0.012%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

37	16.080	2203	2209	2215	rVB5	5834	11605	0.66%	0.063%
38	16.210	2229	2231	2234	rBV4	1763	2275	0.13%	0.012%
39	16.245	2234	2237	2238	rBV3	2903	2424	0.14%	0.013%
40	16.416	2262	2266	2267	rBV4	2877	3338	0.19%	0.018%
41	16.439	2267	2270	2271	rVV3	2322	2295	0.13%	0.012%
42	16.545	2285	2288	2289	rBV3	2418	2359	0.13%	0.013%
43	16.675	2306	2310	2312	rBV5	2969	4280	0.24%	0.023%
44	16.951	2353	2357	2362	rVB7	10606	17272	0.98%	0.093%
45	17.133	2381	2388	2398	rBV	801585	1160070	65.78%	6.262%
46	17.233	2399	2405	2417	rVV	942111	1425232	80.82%	7.693%
47	17.386	2429	2431	2435	rVB5	6195	5051	0.29%	0.027%
48	18.086	2546	2550	2561	rBV2	66524	103336	5.86%	0.558%
49	19.186	2733	2737	2740	rBV3	15255	21515	1.22%	0.116%
50	19.445	2778	2781	2790	rBV3	28311	39402	2.23%	0.213%
51	19.639	2808	2814	2825	rBV2	1086946	1696437	96.20%	9.157%
52	21.315	3095	3099	3106	rVB	124830	183275	10.39%	0.989%
53	21.633	3147	3153	3163	rBV2	700542	1292471	73.29%	6.976%
54	24.768	3677	3686	3699	rVB2	684486	1763501	100.00%	9.519%
55	25.004	3716	3726	3737	rVB	391628	1319908	74.85%	7.124%

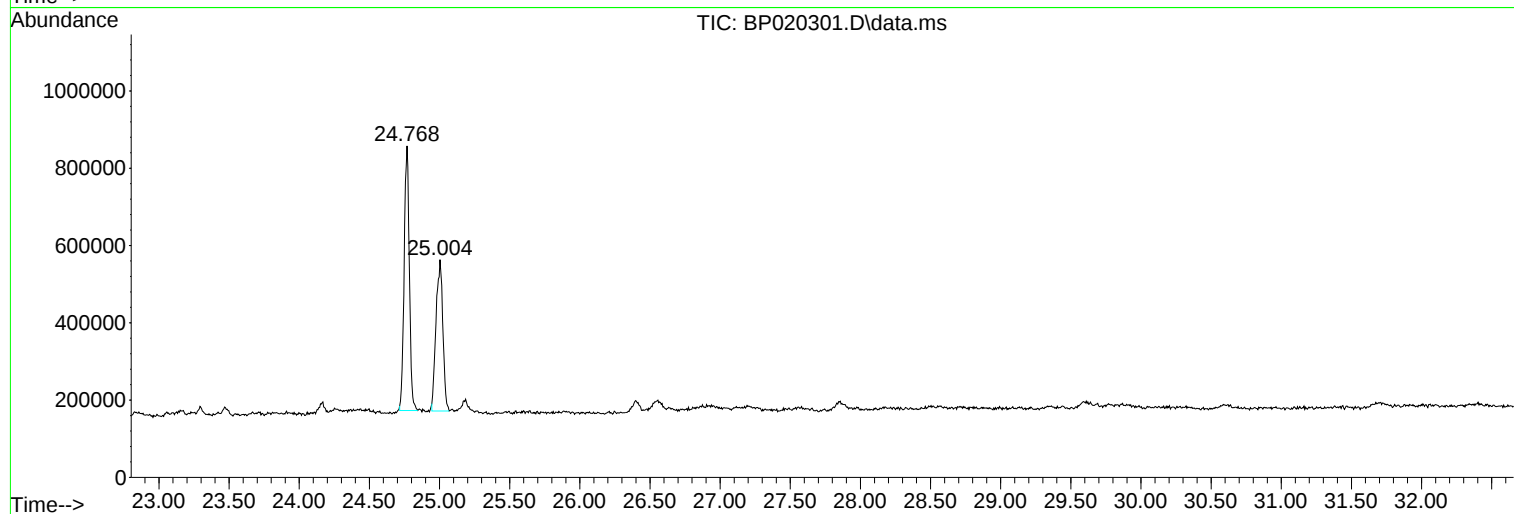
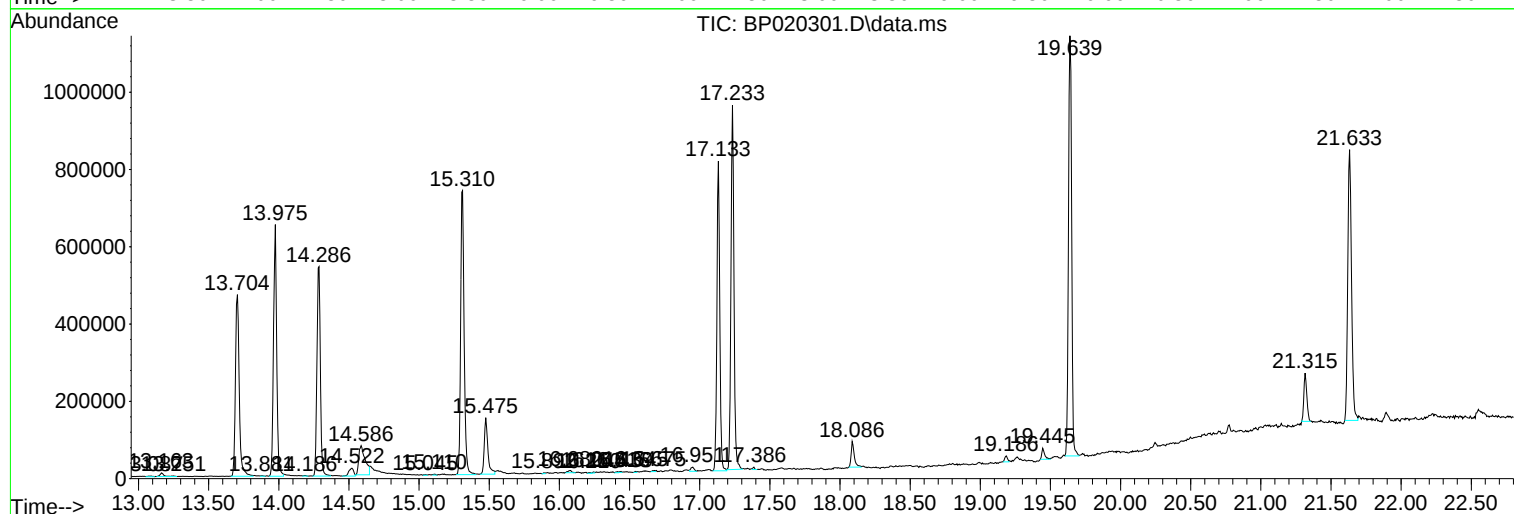
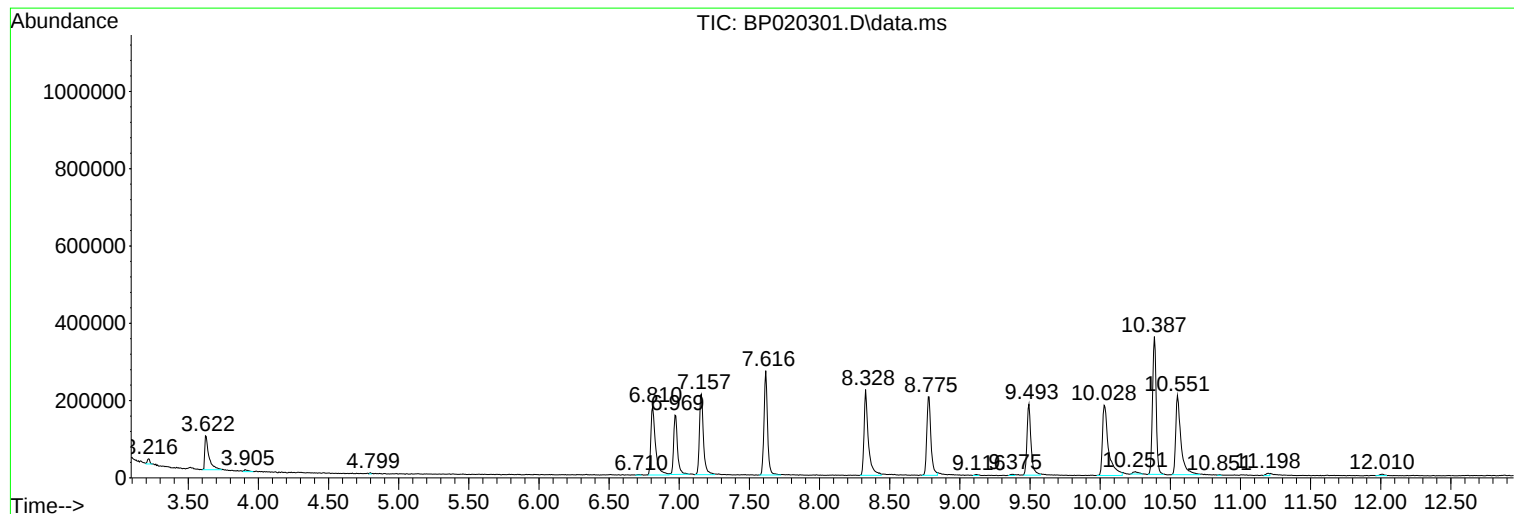
Sum of corrected areas: 18526728

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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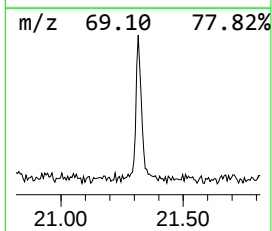
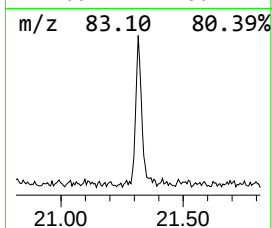
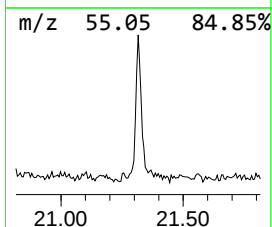
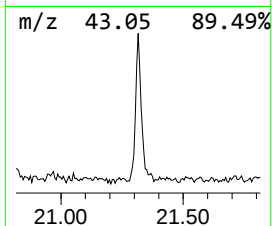
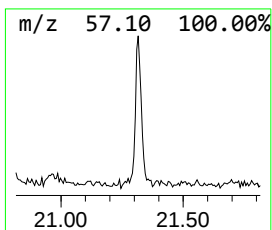
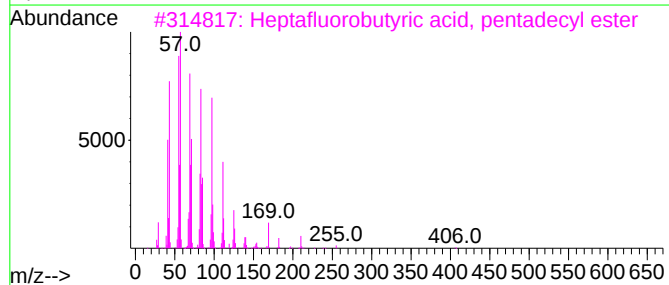
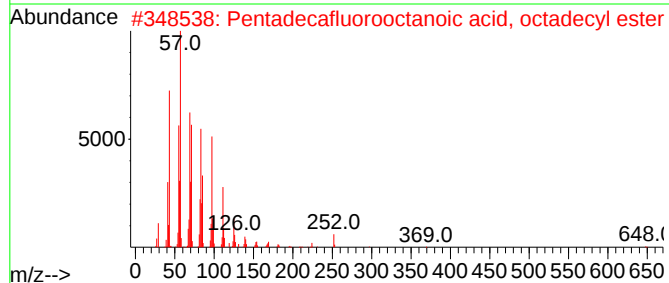
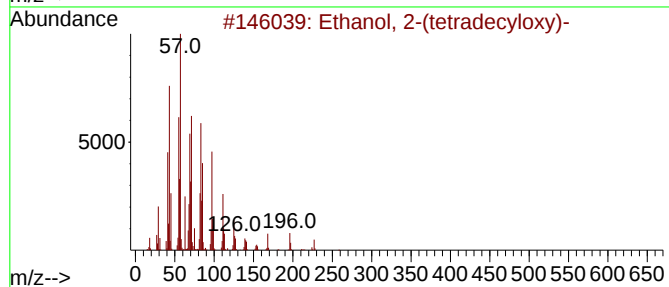
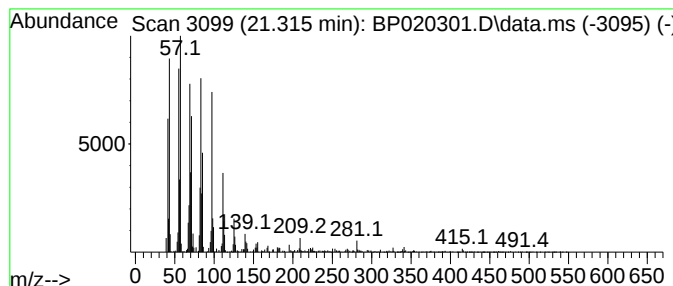
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(tetradecyloxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.316	2.84 ng/ul	183275	Chrysene-d12	21.633

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	95
2			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
4			17-Pentatriacontene	491	C35H70	006971-40-0	91
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(tet...	21.316	2.8	ng/ul	183275	5	21.633	1292470	20.0