

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011512.D
 Acq On : 19 Aug 2022 17:47
 Operator : CG/JU
 Sample : N4221-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KP9

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-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011512.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.069	11	14	19	rVB	23825	25879	2.08%	0.152%
2	3.352	59	62	73	rVB	18455	30798	2.48%	0.181%
3	3.487	82	85	97	rBV	108420	165720	13.34%	0.976%
4	3.693	115	120	125	rBV2	7285	11872	0.96%	0.070%
5	3.781	132	135	139	rVB2	8157	10800	0.87%	0.064%
6	4.216	205	209	216	rVB7	2735	4986	0.40%	0.029%
7	4.493	255	256	262	rBV6	1618	2730	0.22%	0.016%
8	5.011	342	344	346	rVB3	2152	1314	0.11%	0.008%
9	5.516	428	430	435	rBV6	1574	1541	0.12%	0.009%
10	5.681	456	458	461	rVB4	1505	1270	0.10%	0.007%
11	5.805	476	479	480	rBV4	1906	1894	0.15%	0.011%
12	5.940	500	502	506	rVB5	1240	1686	0.14%	0.010%
13	6.246	553	554	558	rVB4	1300	1328	0.11%	0.008%
14	6.740	632	638	654	rVV	292887	468624	37.72%	2.760%
15	6.905	660	666	677	rBV	226883	347107	27.94%	2.044%
16	6.999	679	682	686	rVV	3848	5776	0.46%	0.034%
17	7.087	690	697	710	rBV	309323	481595	38.76%	2.836%
18	7.169	710	711	716	rVB5	2139	1780	0.14%	0.010%
19	7.381	744	747	750	rVB4	1425	1419	0.11%	0.008%
20	7.557	769	777	785	rBV	373127	578775	46.59%	3.409%
21	7.634	786	790	795	rVB6	3402	6070	0.49%	0.036%
22	7.752	808	810	813	rBV4	1041	1448	0.12%	0.009%
23	8.275	892	899	913	rBV	338988	556078	44.76%	3.275%
24	8.557	944	947	949	rVB4	1217	1458	0.12%	0.009%
25	8.722	967	975	986	rBV	273801	455109	36.63%	2.680%
26	8.851	992	997	1001	rBV6	3383	6113	0.49%	0.036%
27	8.893	1003	1004	1007	rVB3	2225	1775	0.14%	0.010%
28	9.122	1039	1043	1049	rVB8	3842	7573	0.61%	0.045%
29	9.434	1086	1096	1105	rBV	224225	374174	30.12%	2.204%
30	9.969	1180	1187	1202	rBV	308530	547280	44.05%	3.223%
31	10.122	1211	1213	1217	rVB5	1546	1822	0.15%	0.011%
32	10.251	1233	1235	1238	rVB4	1442	1571	0.13%	0.009%
33	10.340	1242	1250	1258	rVV	470145	779207	62.72%	4.589%
34	10.498	1266	1277	1291	rBV	289536	517889	41.69%	3.050%
35	10.845	1331	1336	1343	rVB	2953	6180	0.50%	0.036%
36	11.181	1390	1393	1399	rVV6	1305	2828	0.23%	0.017%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
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37	11.281	1405	1410	1411	rBV5	1476	2076	0.17%	0.012%
38	11.298	1411	1413	1416	rVB4	1282	1596	0.13%	0.009%
39	11.487	1444	1445	1448	rVB3	1652	1377	0.11%	0.008%
40	11.510	1448	1449	1453	rBV4	1294	1517	0.12%	0.009%

41	11.557	1453	1457	1460	rVV6	1853	2381	0.19%	0.014%
42	11.616	1463	1467	1471	rVB7	1232	2584	0.21%	0.015%
43	11.945	1519	1523	1530	rVV4	5833	10222	0.82%	0.060%
44	12.063	1542	1543	1548	rVB5	900	1284	0.10%	0.008%
45	12.128	1552	1554	1558	rBV5	1111	1383	0.11%	0.008%

46	12.175	1558	1562	1566	rVB7	1091	1725	0.14%	0.010%
47	12.534	1621	1623	1625	rBV2	1526	1345	0.11%	0.008%
48	12.575	1625	1630	1634	rVV6	1672	2375	0.19%	0.014%
49	13.051	1707	1711	1716	rVB7	2300	4082	0.33%	0.024%
50	13.175	1729	1732	1733	rBV3	1751	1702	0.14%	0.010%

51	13.192	1733	1735	1741	rVB7	2006	3770	0.30%	0.022%
52	13.463	1776	1781	1782	rBV5	1483	2062	0.17%	0.012%
53	13.651	1806	1813	1825	rBV	536056	733360	59.03%	4.319%
54	13.916	1850	1858	1868	rBV	572626	846108	68.10%	4.983%
55	14.016	1874	1875	1880	rVB5	2032	1447	0.12%	0.009%

56	14.139	1890	1896	1899	rBV8	2289	3062	0.25%	0.018%
57	14.228	1904	1911	1920	rVV	646313	905771	72.91%	5.334%
58	14.292	1920	1922	1925	rVB3	2262	1902	0.15%	0.011%
59	14.404	1936	1941	1944	rBV6	2028	3495	0.28%	0.021%
60	14.463	1944	1951	1970	rBV	153934	318727	25.65%	1.877%

61	14.681	1984	1988	1992	rVB7	4312	6576	0.53%	0.039%
62	14.834	2013	2014	2020	rVB6	1180	1445	0.12%	0.009%
63	15.045	2045	2050	2054	rBV8	2476	3555	0.29%	0.021%
64	15.234	2074	2082	2091	rBV	702667	1029314	82.85%	6.062%
65	15.369	2100	2105	2115	rBV2	173799	242166	19.49%	1.426%

66	15.475	2119	2123	2129	rVB9	4432	7132	0.57%	0.042%
67	15.610	2143	2146	2148	rBV4	1883	1909	0.15%	0.011%
68	15.804	2174	2179	2181	rBV6	2123	3767	0.30%	0.022%
69	15.981	2205	2209	2213	rVB7	1863	2728	0.22%	0.016%
70	16.145	2232	2237	2242	rBV9	1882	3316	0.27%	0.020%

71	16.286	2257	2261	2262	rBV4	2088	2266	0.18%	0.013%
72	16.422	2282	2284	2288	rVB4	1778	1620	0.13%	0.010%
73	16.486	2290	2295	2299	rVV8	2830	4257	0.34%	0.025%
74	16.539	2302	2304	2309	rBV5	1033	1552	0.12%	0.009%
75	16.775	2337	2344	2351	rBV5	3491	8081	0.65%	0.048%

76	16.886	2358	2363	2370	rBV5	7630	13506	1.09%	0.080%
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77	16.939	2370	2372	2375	rVB4	2642	2555	0.21%	0.015%
78	16.998	2376	2382	2393	rBV	757173	1026549	82.63%	6.046%
79	17.098	2393	2399	2410	rBV2	802724	1093641	88.03%	6.441%
80	17.216	2415	2419	2426	rVB6	10500	17656	1.42%	0.104%
81	17.569	2474	2479	2480	rBV5	2106	3083	0.25%	0.018%
82	17.698	2497	2501	2506	rVB2	20917	25999	2.09%	0.153%
83	17.916	2534	2538	2548	rBV	46633	71148	5.73%	0.419%
84	17.992	2548	2551	2553	rVV2	3214	4455	0.36%	0.026%
85	18.157	2577	2579	2581	rBV3	2191	1811	0.15%	0.011%
86	18.339	2606	2610	2612	rBV5	3194	4579	0.37%	0.027%
87	18.933	2708	2711	2716	rBV5	11444	17517	1.41%	0.103%
88	19.004	2720	2723	2726	rBV	11183	15079	1.21%	0.089%
89	19.363	2778	2784	2792	rBV	958771	1242367	100.00%	7.317%
90	20.863	3035	3039	3048	rBV	223793	394959	31.79%	2.326%
91	21.127	3079	3084	3089	rBV	925255	1124909	90.55%	6.625%
92	23.268	3442	3448	3455	rBV2	662041	1199459	96.55%	7.064%
93	23.421	3468	3474	3481	rVB	602619	1131889	91.11%	6.666%

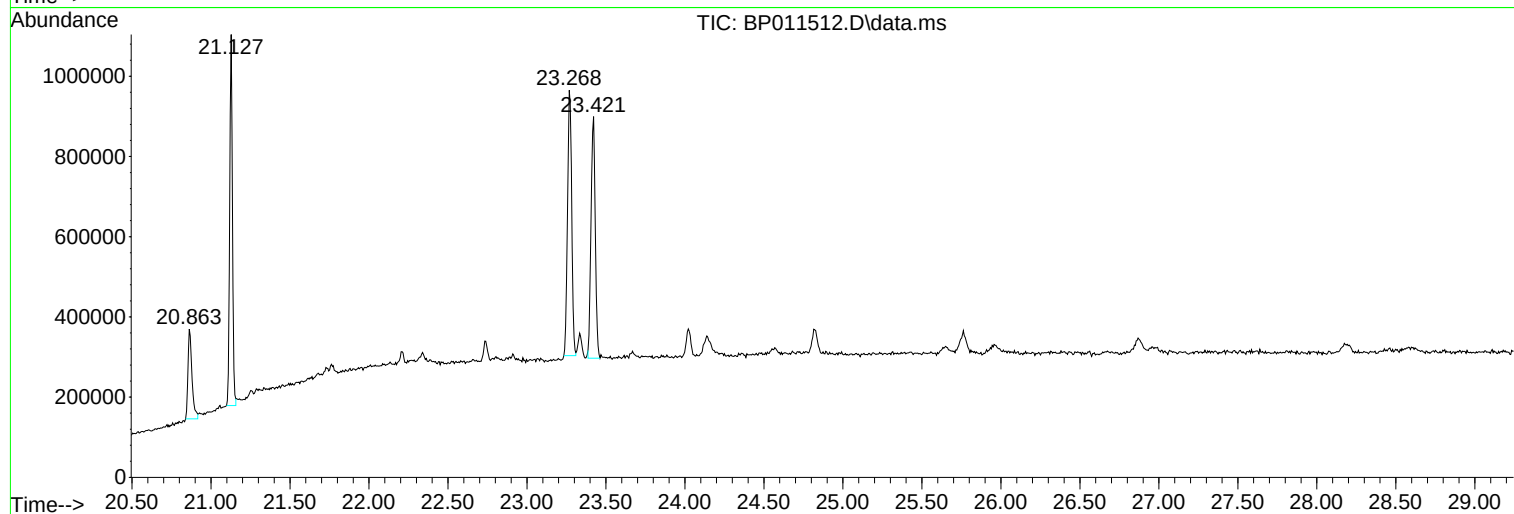
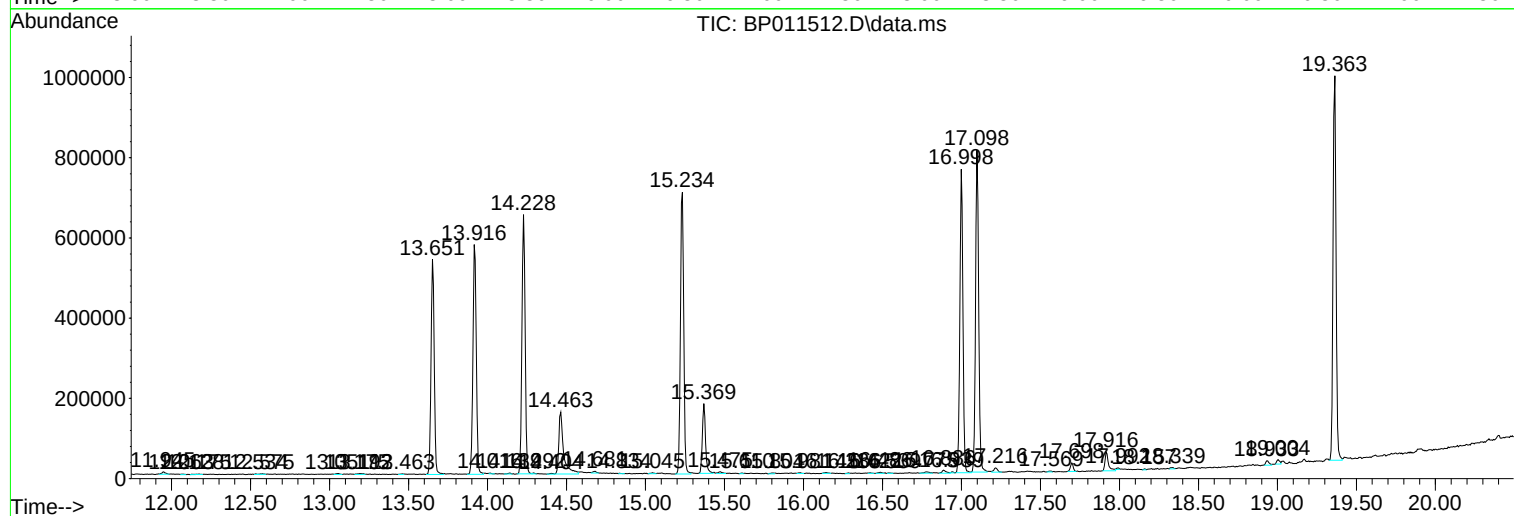
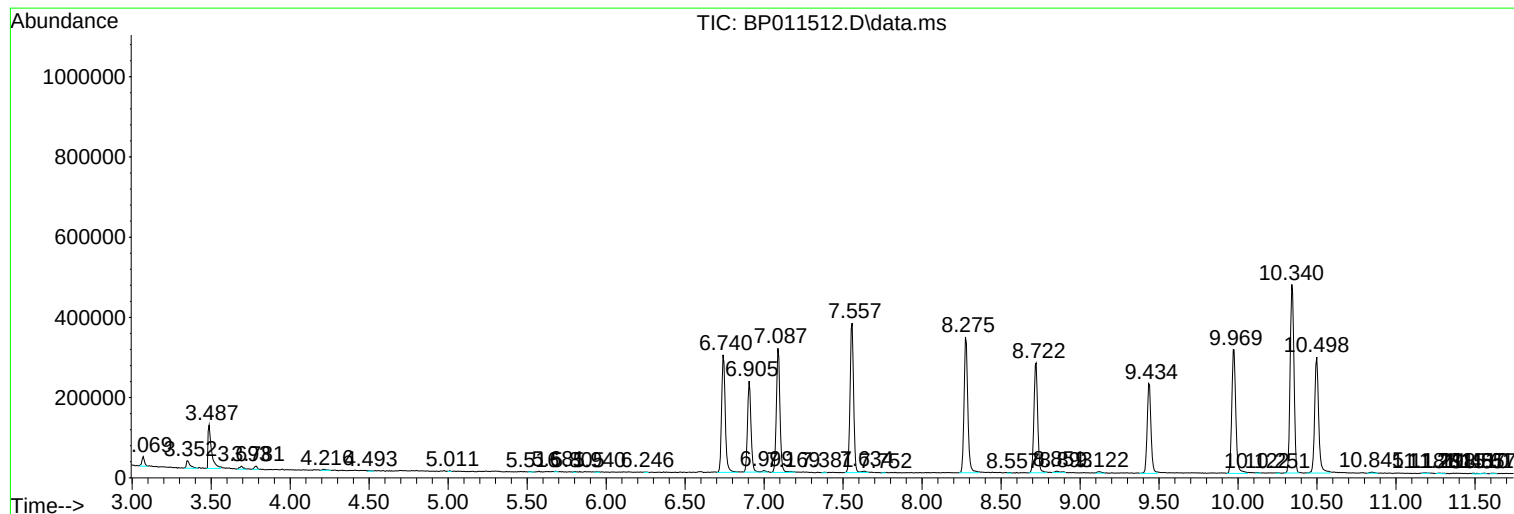
Sum of corrected areas: 16979667

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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



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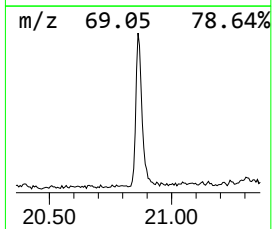
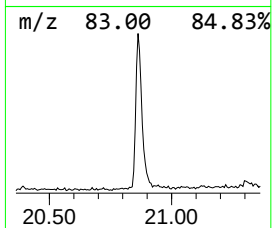
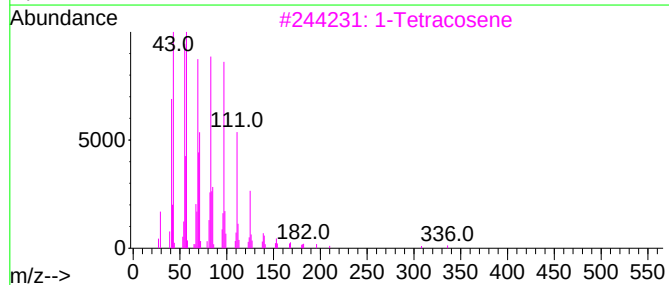
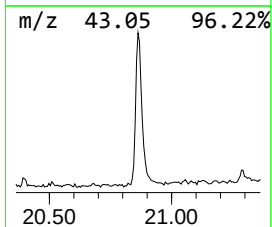
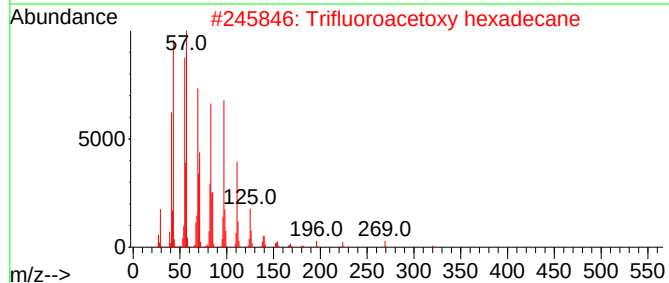
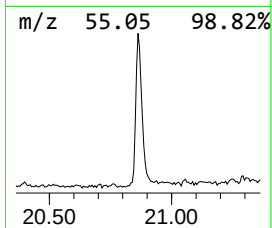
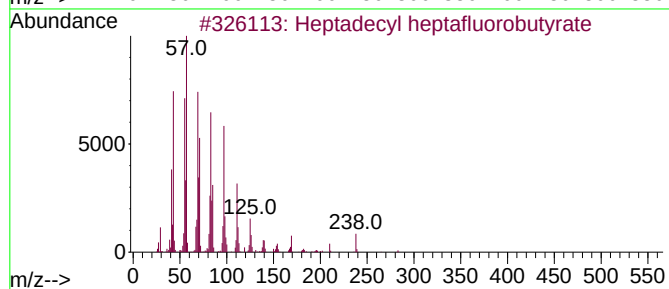
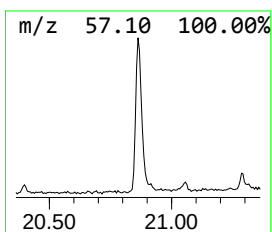
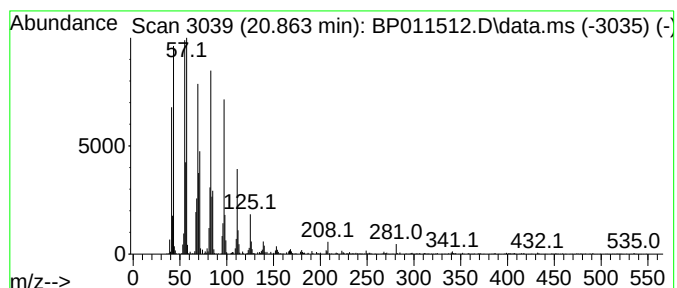
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

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.	
20.863	7.02 ng/ul	394959	Chrysene-d12	21.127	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
3	1-Tetracosene	336	C24H48	010192-32-2	94
4	9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
5	Nonacos-1-ene	406	C29H58	018835-35-3	94



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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
Heptadecyl hept...	20.863	7.0	ng/u1	394959	5	21.127	1124910	20.0