

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072922\
 Data File : BM035992.D
 Acq On : 29 Jul 2022 14:49
 Operator : CG/JU
 Sample : N3790-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBUK0

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_M\Methods\SFAM-EPA-BM072522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM035992.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.269	28	31	42	rVB	90152	121355	2.14%	0.190%
2	3.557	77	80	89	rVB	51067	75684	1.33%	0.119%
3	3.698	101	104	115	rBV	294043	430597	7.58%	0.675%
4	4.028	157	160	166	rVB	22915	30947	0.54%	0.049%
5	5.957	486	488	493	rVB6	5723	7903	0.14%	0.012%
6	6.128	511	517	518	rBV6	3965	5946	0.10%	0.009%
7	7.039	666	672	690	rVB	1152837	1821716	32.07%	2.856%
8	7.210	694	701	710	rVV	987288	1483362	26.11%	2.326%
9	7.286	711	714	720	rVB	11608	18563	0.33%	0.029%
10	7.334	720	722	727	rVB6	4865	6667	0.12%	0.010%
11	7.404	727	734	746	rBV	1202777	1879234	33.08%	2.947%
12	7.492	746	749	754	rVB6	6296	9929	0.17%	0.016%
13	7.875	805	814	823	rBV	998748	1563960	27.53%	2.452%
14	7.945	823	826	831	rVB5	8673	13374	0.24%	0.021%
15	8.586	928	935	949	rBV	1352931	2234184	39.33%	3.503%
16	9.045	1005	1013	1029	rBV	1042287	1721202	30.30%	2.699%
17	9.151	1029	1031	1037	rVB7	4286	5995	0.11%	0.009%
18	9.204	1037	1040	1044	rBV6	8146	10826	0.19%	0.017%
19	9.386	1065	1071	1072	rBV5	3835	6487	0.11%	0.010%
20	9.763	1127	1135	1148	rBV	817337	1373178	24.17%	2.153%
21	10.304	1219	1227	1239	rBV	1277580	2163464	38.09%	3.392%
22	10.680	1284	1291	1300	rBV	1341001	2185790	38.48%	3.427%
23	10.822	1302	1315	1328	rBV	1300543	2200734	38.74%	3.451%
24	11.945	1498	1506	1513	rVB5	10704	26537	0.47%	0.042%
25	12.269	1557	1561	1567	rVB2	23493	36349	0.64%	0.057%
26	12.521	1601	1604	1608	rVB6	4292	6168	0.11%	0.010%
27	12.880	1660	1665	1668	rBV6	10969	18921	0.33%	0.030%
28	13.021	1687	1689	1695	rBV6	4708	7326	0.13%	0.011%
29	13.127	1702	1707	1712	rBV6	8564	10028	0.18%	0.016%
30	13.345	1741	1744	1748	rVB6	5228	7641	0.13%	0.012%
31	13.486	1764	1768	1774	rVB8	8368	13786	0.24%	0.022%
32	13.804	1817	1822	1824	rBV6	4856	6288	0.11%	0.010%
33	13.833	1824	1827	1833	rVB7	5192	10200	0.18%	0.016%
34	13.927	1837	1843	1856	rBV	2229993	3033647	53.41%	4.757%
35	14.210	1884	1891	1899	rVV	2608577	3792601	66.77%	5.947%
36	14.286	1902	1904	1910	rVB7	8436	12975	0.23%	0.020%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
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37	14.415	1923	1926	1930	rVB6	8049	8722	0.15%	0.014%
38	14.468	1930	1935	1936	rBV5	4347	5912	0.10%	0.009%
39	14.515	1936	1943	1953	rBV	1873691	2747688	48.37%	4.308%
40	14.710	1970	1976	1999	rVV	840300	1347493	23.72%	2.113%
41	14.933	2011	2014	2019	rVB6	13710	20338	0.36%	0.032%
42	15.309	2073	2078	2079	rBV4	11102	13088	0.23%	0.021%
43	15.357	2084	2086	2091	rVB6	6347	8514	0.15%	0.013%
44	15.504	2103	2111	2122	rBV	3409819	4787594	84.28%	7.507%
45	15.621	2125	2131	2147	rBV	777614	1091717	19.22%	1.712%
46	15.745	2147	2152	2156	rBV	39300	60961	1.07%	0.096%
47	15.874	2169	2174	2178	rBV7	4850	8016	0.14%	0.013%
48	16.068	2204	2207	2209	rBV4	5069	6277	0.11%	0.010%
49	16.186	2223	2227	2230	rBV6	5363	8650	0.15%	0.014%
50	16.221	2230	2233	2236	rVV5	12177	18089	0.32%	0.028%
51	16.286	2240	2244	2249	rVB6	14044	21826	0.38%	0.034%
52	16.398	2257	2263	2268	rBV10	9523	18320	0.32%	0.029%
53	16.562	2286	2291	2295	rBV7	11436	18505	0.33%	0.029%
54	16.733	2316	2320	2324	rBV2	30233	38756	0.68%	0.061%
55	16.933	2350	2354	2360	rBV7	8064	14987	0.26%	0.023%
56	17.015	2364	2368	2370	rBV4	15825	22448	0.40%	0.035%
57	17.098	2380	2382	2385	rBV4	5960	6759	0.12%	0.011%
58	17.174	2393	2395	2399	rVB5	6823	8952	0.16%	0.014%
59	17.251	2402	2408	2415	rBV	2240787	3147568	55.41%	4.935%
60	17.351	2419	2425	2437	rVV2	3534733	5013439	88.26%	7.861%
61	17.445	2437	2441	2446	rVB8	22674	32319	0.57%	0.051%
62	17.633	2471	2473	2479	rVB7	8548	12427	0.22%	0.019%
63	17.803	2498	2502	2509	rVB	24991	36196	0.64%	0.057%
64	17.927	2520	2523	2528	rVB3	26717	32047	0.56%	0.050%
65	18.150	2557	2561	2565	rBV3	36332	47884	0.84%	0.075%
66	18.221	2571	2573	2577	rVB	10361	13663	0.24%	0.021%
67	18.750	2660	2663	2667	rVB2	25876	30033	0.53%	0.047%
68	19.209	2736	2741	2744	rBV2	46229	63055	1.11%	0.099%
69	19.280	2749	2753	2756	rBV	43630	55505	0.98%	0.087%
70	19.544	2796	2798	2801	rBV3	24147	23967	0.42%	0.038%
71	19.645	2809	2815	2821	rBV	4181028	5680330	100.00%	8.907%
72	19.692	2821	2823	2829	rVB2	38242	49775	0.88%	0.078%
73	20.644	2982	2985	2989	rBV	136244	149210	2.63%	0.234%
74	21.150	3067	3071	3077	rBV2	295439	398013	7.01%	0.624%
75	21.433	3114	3119	3125	rBV	2649876	3406303	59.97%	5.341%
76	23.650	3489	3496	3503	rBV	2983885	5573313	98.12%	8.739%

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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

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Peak separation: 5

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Title : SVOA CALIBRATION

77 23.803 3515 3522 3530 rVB2 1674380 3372897 59.38% 5.289%

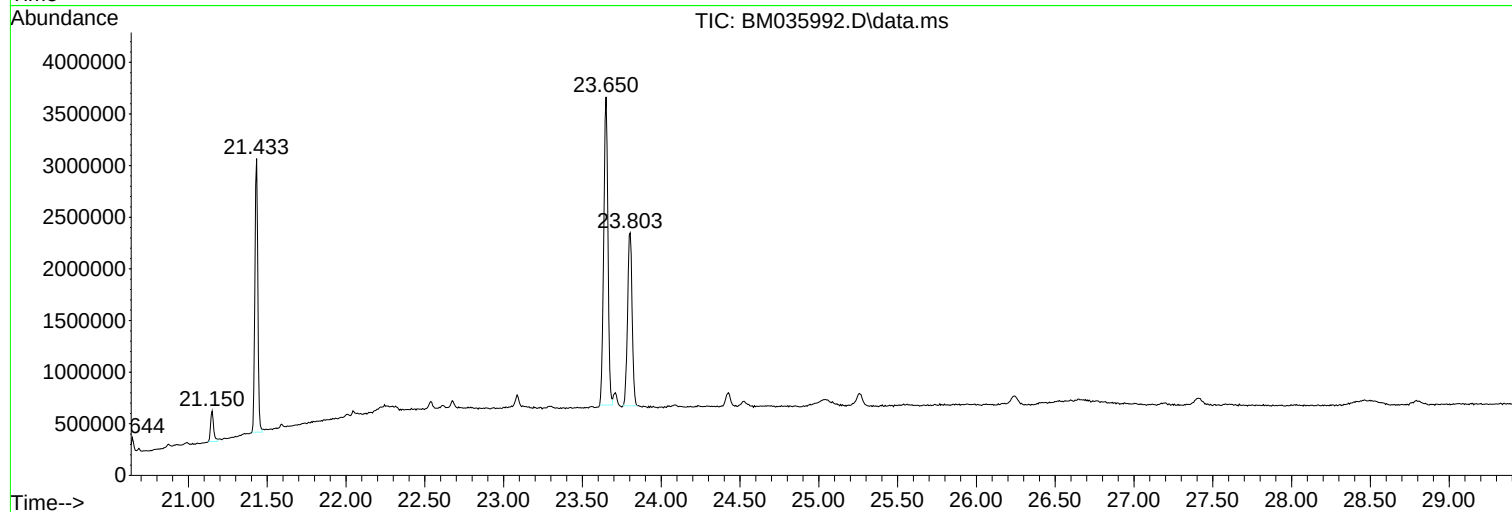
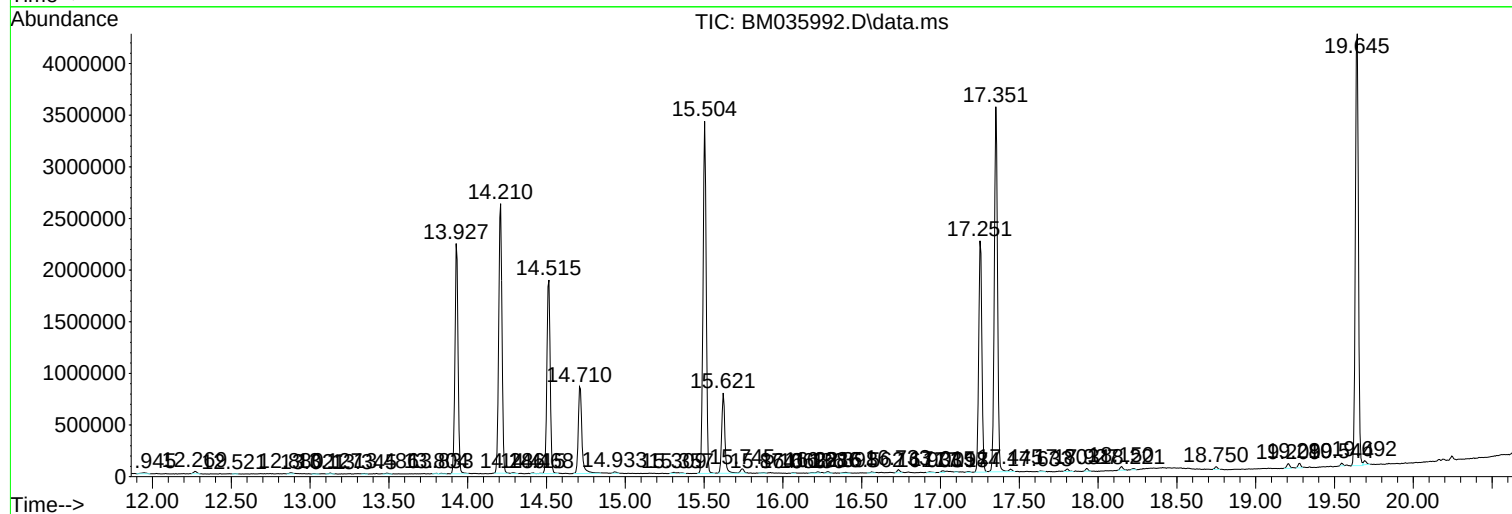
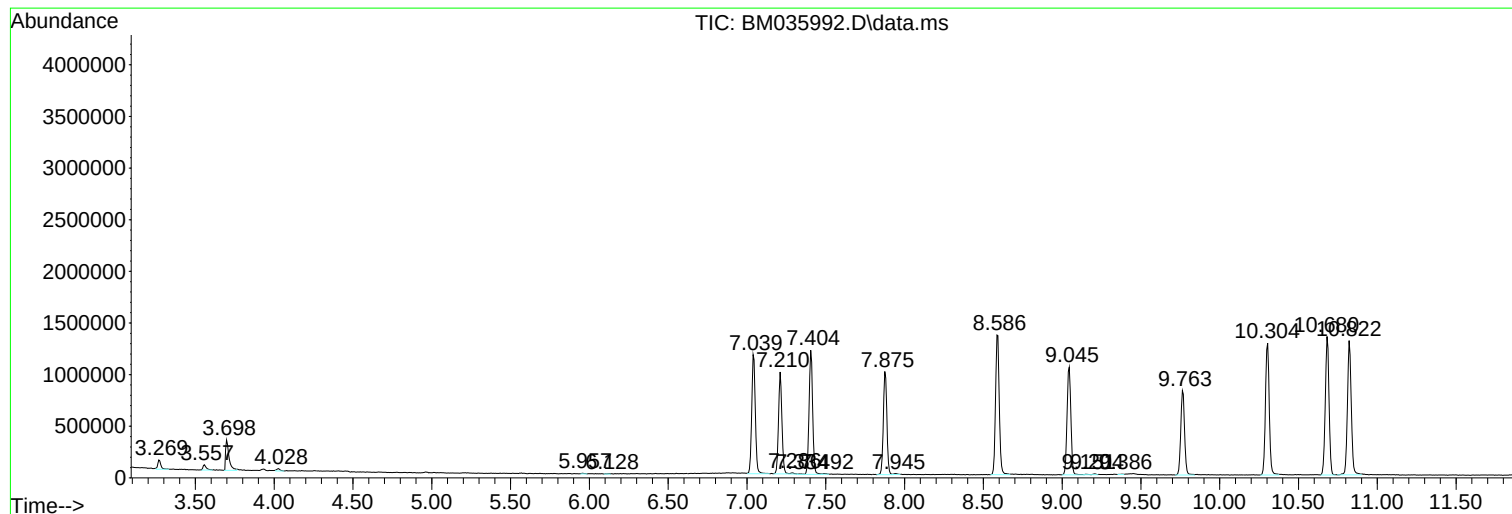
Sum of corrected areas: 63775120

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Instrument :
BNA_M
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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



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BNA_M
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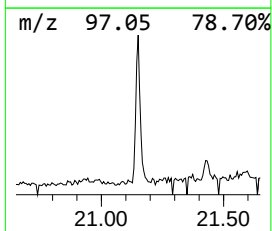
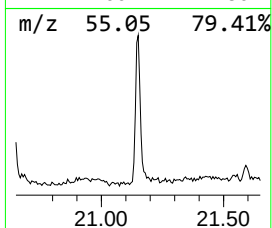
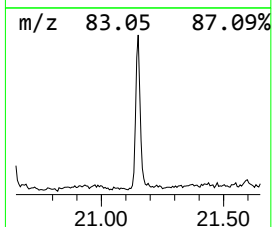
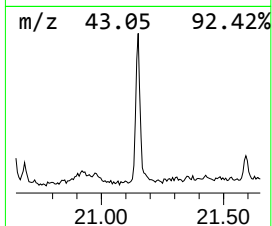
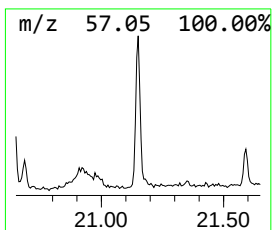
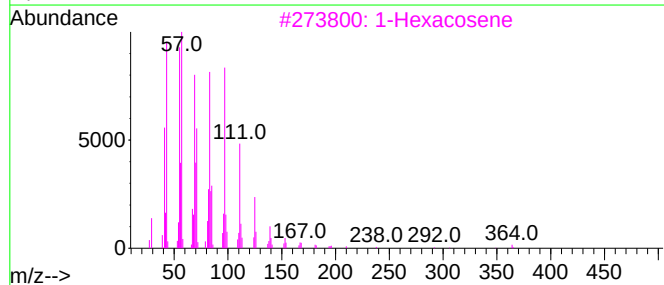
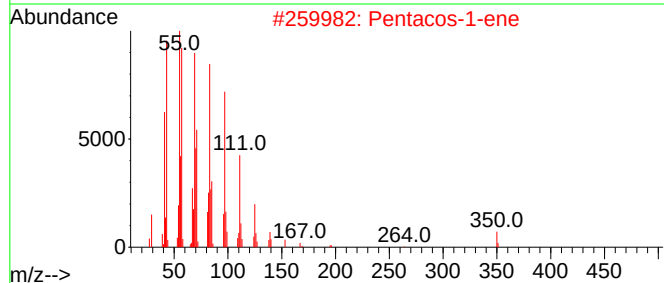
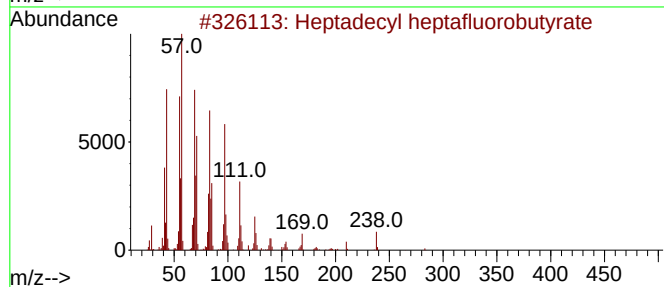
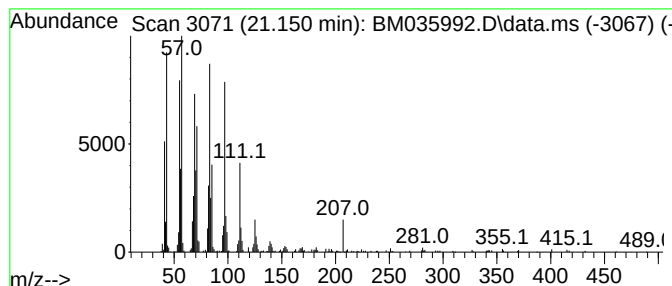
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
Quant Title : SVOA CALIBRATION

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.
21.150	2.34 ng/ul	398013	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			Pentacos-1-ene	350	C25H50	016980-85-1	91
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	90
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	90



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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...	21.150	2.3	ng/u1	398013	5	21.433	3406300	20.0