

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012882.D
 Acq On : 30 Nov 2022 16:39
 Operator : CG/JU
 Sample : N5725-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB55

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-BP112522.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012882.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	31	35	46	rVB	115033	177755	1.12%	0.124%
2	3.828	106	109	121	rBV	271788	435183	2.74%	0.304%
3	4.058	143	148	158	rVB	62191	104014	0.66%	0.073%
4	4.158	160	165	170	rVB	105044	147284	0.93%	0.103%
5	4.670	247	252	260	rVB	56642	97944	0.62%	0.068%
6	5.628	411	415	421	rBV3	15705	30956	0.19%	0.022%
7	6.899	627	631	635	rBV4	12053	18923	0.12%	0.013%
8	7.158	669	675	684	rVB	458969	697491	4.39%	0.488%
9	7.334	698	705	715	rBV	1997605	3046780	19.19%	2.130%
10	7.534	732	739	749	rBV	1607297	2568890	16.18%	1.796%
11	7.628	751	755	762	rVB9	14351	27167	0.17%	0.019%
12	8.011	813	820	826	rBV	2205842	3439235	21.66%	2.404%
13	8.069	826	830	838	rVB	314225	461048	2.90%	0.322%
14	8.228	851	857	865	rBV	10476	27794	0.18%	0.019%
15	8.387	878	884	889	rBV6	9662	18619	0.12%	0.013%
16	8.710	932	939	952	rVV	1381617	2122641	13.37%	1.484%
17	8.834	956	960	968	rVB	35749	60802	0.38%	0.042%
18	9.110	1000	1007	1011	rBV	38677	68968	0.43%	0.048%
19	9.175	1011	1018	1028	rVV	2290031	3575213	22.52%	2.499%
20	9.269	1028	1034	1041	rVV6	32695	74237	0.47%	0.052%
21	9.334	1041	1045	1050	rVB7	10194	20181	0.13%	0.014%
22	9.593	1084	1089	1093	rVB5	16817	31257	0.20%	0.022%
23	9.734	1111	1113	1120	rVB7	9758	17643	0.11%	0.012%
24	9.899	1133	1141	1152	rBV	1491841	2457413	15.48%	1.718%
25	10.122	1175	1179	1185	rVB9	8182	16534	0.10%	0.012%
26	10.293	1203	1208	1211	rBV6	8224	16168	0.10%	0.011%
27	10.440	1222	1233	1245	rBV	2524882	4183911	26.35%	2.924%
28	10.599	1256	1260	1268	rBV9	11464	20276	0.13%	0.014%
29	10.687	1268	1275	1278	rBV8	13696	27436	0.17%	0.019%
30	10.734	1278	1283	1289	rVB6	23499	47210	0.30%	0.033%
31	10.828	1289	1299	1313	rBV	3290105	5706578	35.94%	3.989%
32	10.957	1313	1321	1336	rBV	1494826	2568754	16.18%	1.795%
33	11.481	1405	1410	1416	rVB3	12748	26361	0.17%	0.018%
34	11.593	1424	1429	1433	rBV8	10249	19947	0.13%	0.014%
35	12.075	1506	1511	1518	rBV8	9837	19173	0.12%	0.013%
36	12.404	1555	1567	1573	rBV2	57521	114618	0.72%	0.080%

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Title : SVOA CALIBRATION

37	12.634	1599	1606	1610	rBV3	14779	29372	0.19%	0.021%
38	12.822	1634	1638	1647	rBV10	10684	27880	0.18%	0.019%
39	12.934	1652	1657	1664	rVB	357036	513424	3.23%	0.359%
40	13.257	1706	1712	1718	rVB6	23196	42726	0.27%	0.030%
41	13.481	1746	1750	1754	rBV4	25380	40736	0.26%	0.028%
42	13.651	1768	1779	1786	rVB	365833	537993	3.39%	0.376%
43	13.887	1812	1819	1824	rBV9	11713	24242	0.15%	0.017%
44	13.963	1828	1832	1837	rVB6	15853	21431	0.13%	0.015%
45	14.051	1839	1847	1853	rBV	6015272	8177654	51.51%	5.716%
46	14.222	1872	1876	1881	rVB3	26536	38453	0.24%	0.027%
47	14.340	1889	1896	1908	rVV	6550464	9701576	61.11%	6.781%
48	14.645	1939	1948	1955	rBV	5709242	8260259	52.03%	5.774%
49	14.716	1955	1960	1965	rVB4	50409	93334	0.59%	0.065%
50	14.822	1973	1978	1989	rBV	245105	417356	2.63%	0.292%
51	15.051	2012	2017	2030	rVB4	59085	119622	0.75%	0.084%
52	15.316	2058	2062	2067	rVB2	37568	56230	0.35%	0.039%
53	15.440	2079	2083	2084	rBV2	21191	22763	0.14%	0.016%
54	15.640	2110	2117	2128	rBV	9337459	13207106	83.19%	9.231%
55	15.745	2129	2135	2151	rVV	1553428	2150454	13.55%	1.503%
56	15.881	2153	2158	2164	rVV	44645	80986	0.51%	0.057%
57	15.939	2164	2168	2173	rVV2	31985	48588	0.31%	0.034%
58	15.998	2173	2178	2188	rVB	52664	85403	0.54%	0.060%
59	16.210	2210	2214	2218	rBV6	16224	26315	0.17%	0.018%
60	16.992	2342	2347	2352	rBV2	17404	30822	0.19%	0.022%
61	17.075	2357	2361	2369	rVB9	13841	26513	0.17%	0.019%
62	17.186	2376	2380	2387	rVB3	13893	27463	0.17%	0.019%
63	17.357	2404	2409	2410	rBV5	13040	19312	0.12%	0.013%
64	17.398	2410	2416	2427	rVV	6320112	9551678	60.16%	6.676%
65	17.498	2427	2433	2444	rVV	10031156	14587703	91.88%	10.196%
66	17.586	2444	2448	2453	rVB2	70341	104295	0.66%	0.073%
67	17.681	2460	2464	2471	rVB	42304	55029	0.35%	0.038%
68	17.804	2482	2485	2490	rVB	15592	22126	0.14%	0.015%
69	18.057	2523	2528	2537	rBV	493597	661236	4.16%	0.462%
70	18.269	2560	2564	2568	rBV7	21858	31805	0.20%	0.022%
71	18.345	2574	2577	2583	rVB	72575	91910	0.58%	0.064%
72	18.516	2604	2606	2611	rVB5	12721	15967	0.10%	0.011%
73	19.304	2736	2740	2746	rBV	117789	167171	1.05%	0.117%
74	19.375	2747	2752	2756	rVV	108577	162242	1.02%	0.113%
75	19.416	2757	2759	2763	rVB3	31411	33301	0.21%	0.023%
76	19.645	2795	2798	2803	rVB4	51059	59261	0.37%	0.041%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Title : SVOA CALIBRATION

77	19.728	2806	2812	2824	rBV	11925869	15876312	100.00%	11.097%
78	21.486	3105	3111	3121	rBV	5969973	8033364	50.60%	5.615%
79	23.833	3503	3510	3525	rBV2	5255064	10647160	67.06%	7.442%
80	23.998	3531	3538	3553	rVB	3104397	6646574	41.86%	4.646%

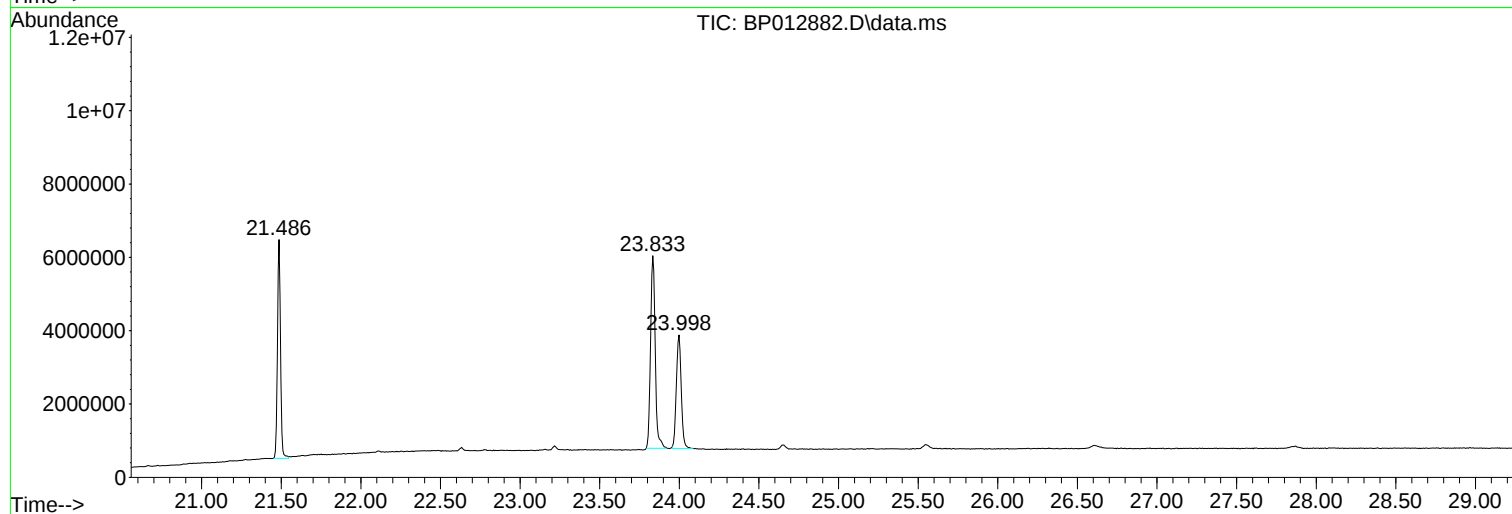
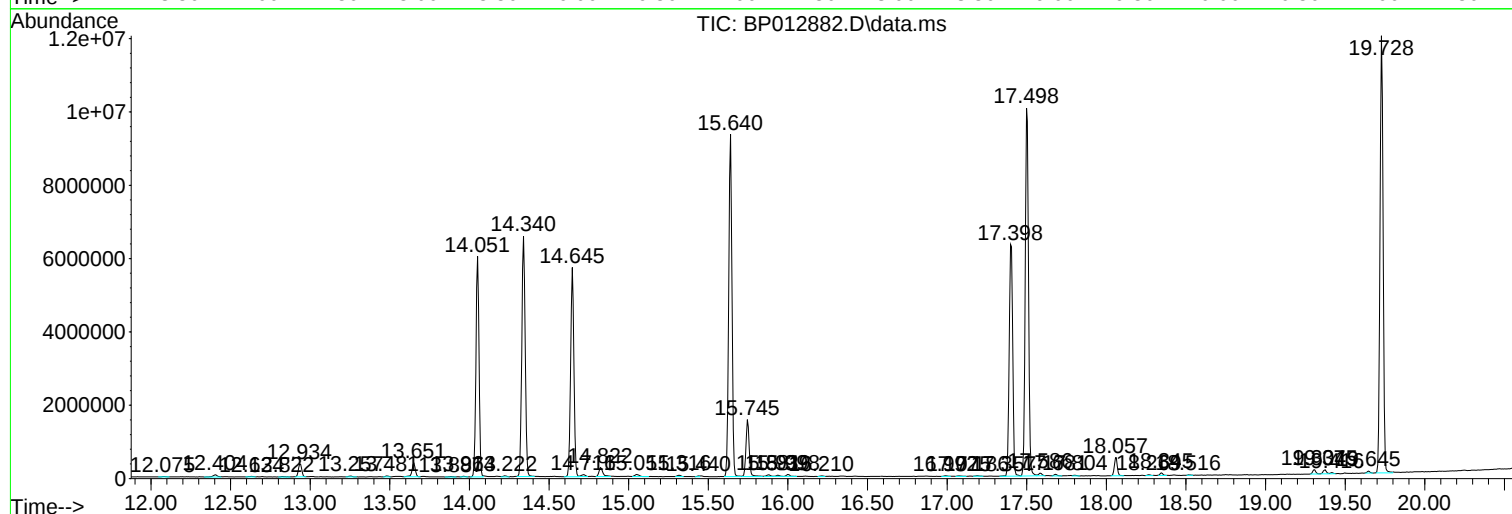
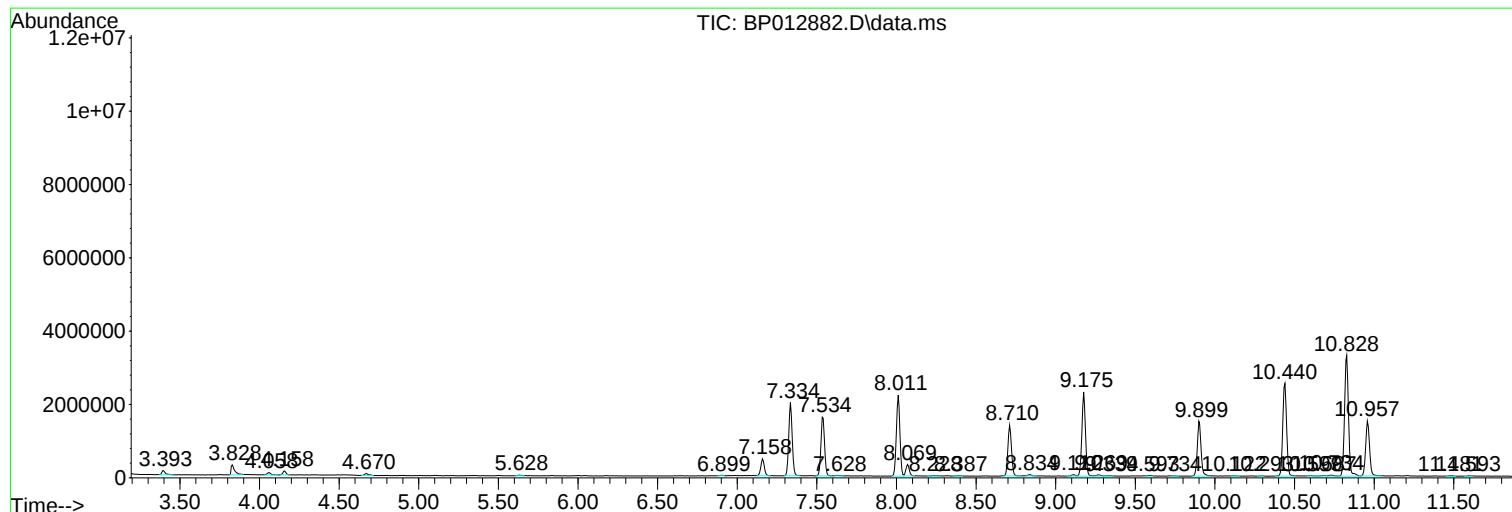
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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



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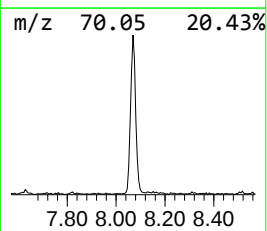
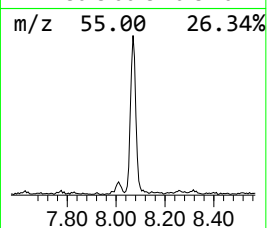
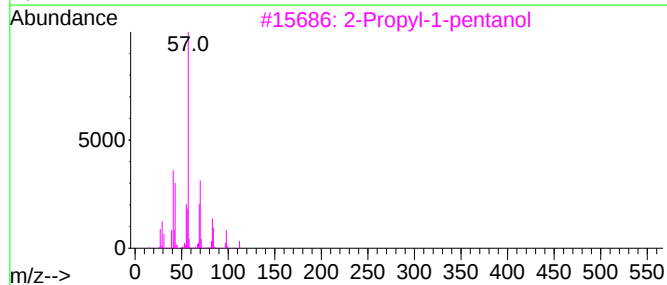
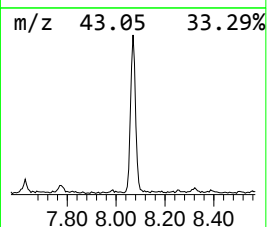
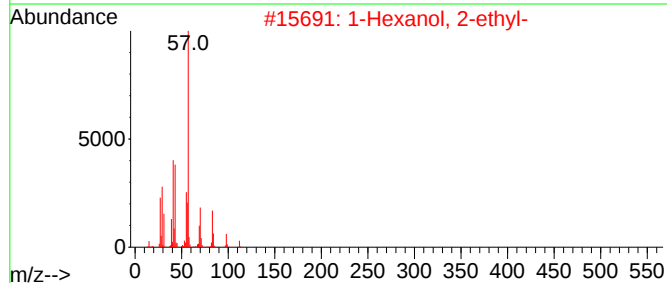
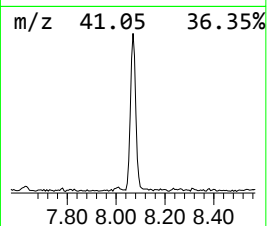
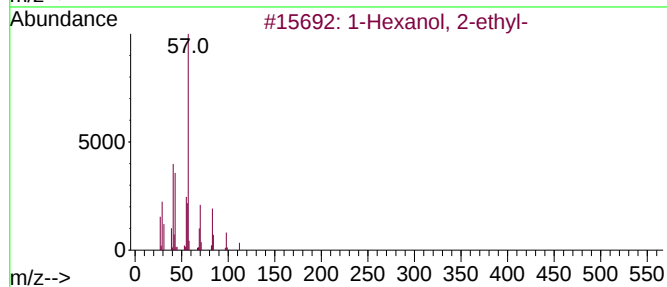
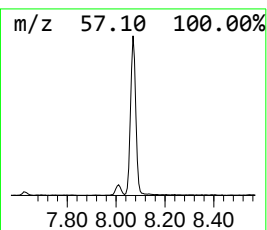
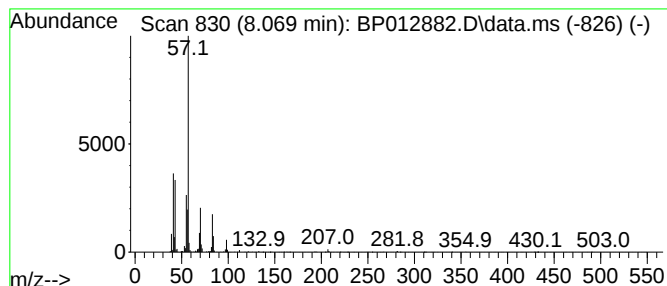
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.069	2.68 ng/ul	461048	1,4-Dichlorobenzene-d4	8.011

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
3	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	59



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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
1-Hexanol, 2-et...	8.069	2.7	ng/ul	461048	1	8.011	3439240	20.0