

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012892.D
 Acq On : 30 Nov 2022 22:56
 Operator : CG/JU
 Sample : N5759-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB80

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: BP012892.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	44	rVB	116801	163937	0.98%	0.110%
2	3.822	106	108	122	rBV	264748	458340	2.75%	0.308%
3	4.057	143	148	154	rVB	59740	92810	0.56%	0.062%
4	4.152	160	164	169	rVB	58270	79922	0.48%	0.054%
5	4.669	248	252	260	rVB	34725	59668	0.36%	0.040%
6	5.099	320	325	330	rBV5	10310	19774	0.12%	0.013%
7	5.175	334	338	345	rVB10	14095	21284	0.13%	0.014%
8	5.634	410	416	417	rBV2	31663	46627	0.28%	0.031%
9	6.004	473	479	483	rBV4	13057	21246	0.13%	0.014%
10	6.646	583	588	596	rVB5	20192	34536	0.21%	0.023%
11	6.740	596	604	613	rBV4	28056	73775	0.44%	0.050%
12	7.087	659	663	668	rBV3	17013	27771	0.17%	0.019%
13	7.157	668	675	683	rVB	523820	802305	4.82%	0.540%
14	7.334	698	705	714	rBV	1960925	3048195	18.31%	2.050%
15	7.534	731	739	747	rVV	1755591	2764263	16.60%	1.859%
16	7.651	751	759	764	rVB2	39247	76422	0.46%	0.051%
17	8.010	811	820	826	rBV	1917549	3063478	18.40%	2.060%
18	8.069	826	830	836	rVV	276323	410899	2.47%	0.276%
19	8.122	836	839	844	rVB4	10954	17697	0.11%	0.012%
20	8.663	920	931	932	rBV10	13259	28427	0.17%	0.019%
21	8.710	932	939	949	rVB	1544259	2398746	14.41%	1.613%
22	8.834	955	960	968	rVB	39334	69067	0.41%	0.046%
23	9.110	1000	1007	1011	rBV3	61626	115434	0.69%	0.078%
24	9.175	1011	1018	1027	rVB	2201131	3546990	21.30%	2.386%
25	9.898	1132	1141	1152	rBV	1665813	2700378	16.22%	1.816%
26	10.228	1190	1197	1202	rBV10	9540	18302	0.11%	0.012%
27	10.434	1223	1232	1252	rVB	2609960	4487781	26.95%	3.018%
28	10.598	1254	1260	1266	rBV9	14202	27165	0.16%	0.018%
29	10.686	1269	1275	1278	rBV6	17518	30000	0.18%	0.020%
30	10.728	1278	1282	1290	rVB7	26511	50710	0.30%	0.034%
31	10.822	1290	1298	1305	rBV	2896898	4981695	29.92%	3.351%
32	10.881	1305	1308	1313	rVB3	46031	69291	0.42%	0.047%
33	10.957	1313	1321	1335	rBV	1068454	1823861	10.95%	1.227%
34	11.204	1360	1363	1369	rVB7	15802	24765	0.15%	0.017%
35	11.481	1406	1410	1415	rVB3	12149	20552	0.12%	0.014%
36	11.757	1452	1457	1462	rBV	18461	31182	0.19%	0.021%

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Start Thrs: 0.2

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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

37	12.075	1508	1511	1518	rVB7	9841	19748	0.12%	0.013%
38	12.175	1523	1528	1532	rVV	49481	80162	0.48%	0.054%
39	12.216	1532	1535	1543	rVB9	19758	37177	0.22%	0.025%
40	12.404	1559	1567	1573	rBV3	70609	139062	0.84%	0.094%

41	12.480	1573	1580	1588	rVB7	23679	63551	0.38%	0.043%
42	12.628	1599	1605	1609	rBV4	15168	22650	0.14%	0.015%
43	12.828	1633	1639	1644	rBV8	12142	26878	0.16%	0.018%
44	12.933	1652	1657	1665	rVB	443235	616983	3.71%	0.415%
45	13.010	1665	1670	1674	rVB6	14715	21420	0.13%	0.014%

46	13.263	1704	1713	1721	rVB3	45693	107125	0.64%	0.072%
47	13.451	1742	1745	1746	rVV3	18241	21208	0.13%	0.014%
48	13.480	1746	1750	1755	rVV	94269	141956	0.85%	0.095%
49	13.533	1755	1759	1765	rVV5	21606	38353	0.23%	0.026%
50	13.651	1765	1779	1784	rVV	409888	673297	4.04%	0.453%

51	13.710	1786	1789	1794	rVB5	15041	23137	0.14%	0.016%
52	13.963	1828	1832	1837	rVB8	12344	18805	0.11%	0.013%
53	14.051	1839	1847	1857	rBV	5981340	8115094	48.74%	5.458%
54	14.163	1862	1866	1869	rVV6	16810	27606	0.17%	0.019%
55	14.222	1872	1876	1883	rVB2	27945	43275	0.26%	0.029%

56	14.339	1889	1896	1910	rBV	6781001	9800040	58.86%	6.591%
57	14.504	1920	1924	1927	rBV3	21847	30012	0.18%	0.020%
58	14.645	1939	1948	1956	rBV	5105454	7530480	45.23%	5.065%
59	14.722	1956	1961	1966	rVB6	30854	50053	0.30%	0.034%
60	14.822	1972	1978	1984	rBV	309829	505919	3.04%	0.340%

61	15.057	2013	2018	2024	rVB7	36869	67501	0.41%	0.045%
62	15.316	2054	2062	2068	rBV	49454	79402	0.48%	0.053%
63	15.380	2068	2073	2077	rVV3	19728	30682	0.18%	0.021%
64	15.439	2077	2083	2096	rVB2	34497	75153	0.45%	0.051%
65	15.639	2110	2117	2124	rBV	8859880	12865716	77.27%	8.653%

66	15.745	2129	2135	2150	rVV	2018204	2642934	15.87%	1.778%
67	15.874	2153	2157	2163	rVV2	47274	72922	0.44%	0.049%
68	15.933	2163	2167	2172	rVB4	36652	53779	0.32%	0.036%
69	15.998	2173	2178	2187	rVB	57753	83473	0.50%	0.056%
70	16.210	2211	2214	2219	rVB7	16150	19795	0.12%	0.013%

71	16.868	2322	2326	2332	rVB7	11706	22308	0.13%	0.015%
72	16.992	2341	2347	2353	rBV8	16709	29061	0.17%	0.020%
73	17.051	2353	2357	2369	rVV3	41214	76892	0.46%	0.052%
74	17.180	2376	2379	2384	rBV2	13447	17642	0.11%	0.012%
75	17.398	2410	2416	2427	rBV	6455452	9172043	55.09%	6.169%

76	17.498	2427	2433	2444	rVV2	9996164	14179244	85.16%	9.537%
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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	17.586	2444	2448	2453	rVB4	48636	71416	0.43%	0.048%
78	17.680	2459	2464	2470	rVB	284431	347317	2.09%	0.234%
79	18.057	2523	2528	2534	rVV	532873	650216	3.91%	0.437%
80	18.263	2560	2563	2569	rVB3	53317	63746	0.38%	0.043%
81	18.345	2573	2577	2585	rVB	95215	140862	0.85%	0.095%
82	18.921	2671	2675	2680	rBV	63191	87469	0.53%	0.059%
83	19.304	2736	2740	2744	rBV3	117845	159125	0.96%	0.107%
84	19.368	2747	2751	2756	rVV	114990	161540	0.97%	0.109%
85	19.415	2756	2759	2762	rVB5	27905	35540	0.21%	0.024%
86	19.492	2770	2772	2776	rBV3	35012	43694	0.26%	0.029%
87	19.645	2795	2798	2802	rVB3	56301	65425	0.39%	0.044%
88	19.727	2806	2812	2822	rBV	12739255	16650382	100.00%	11.199%
89	21.486	3105	3111	3121	rBV	7098493	9426022	56.61%	6.340%
90	23.833	3502	3510	3526	rBV	6316096	13470193	80.90%	9.060%
91	23.998	3531	3538	3550	rVB	3786808	7958359	47.80%	5.353%

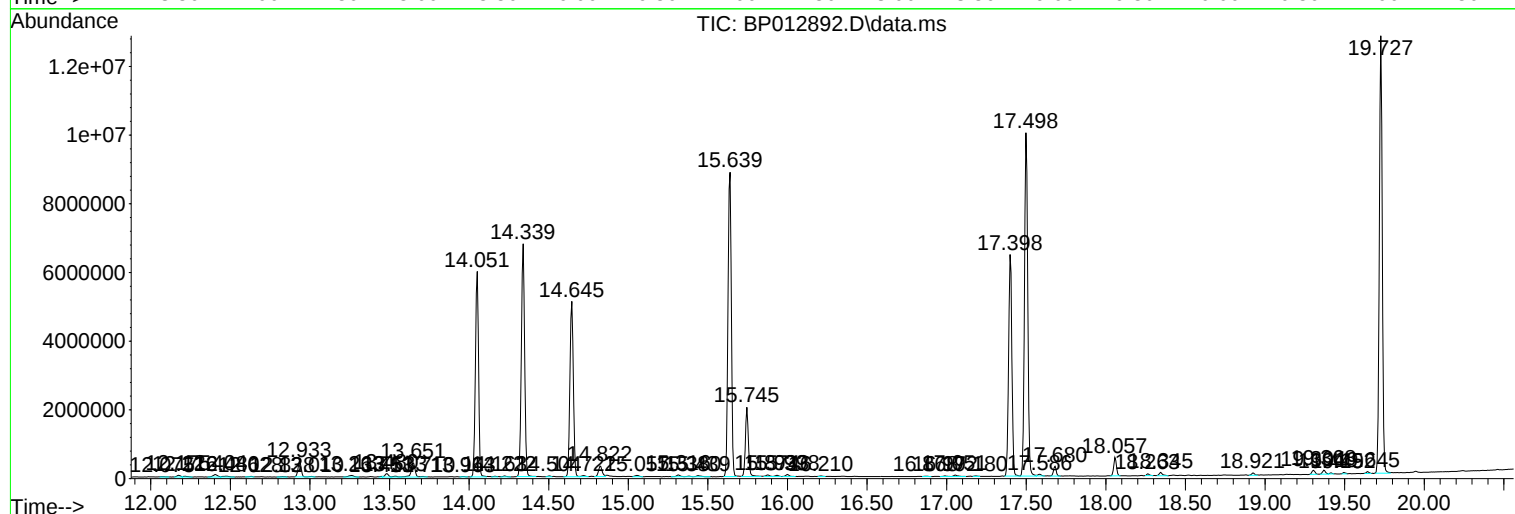
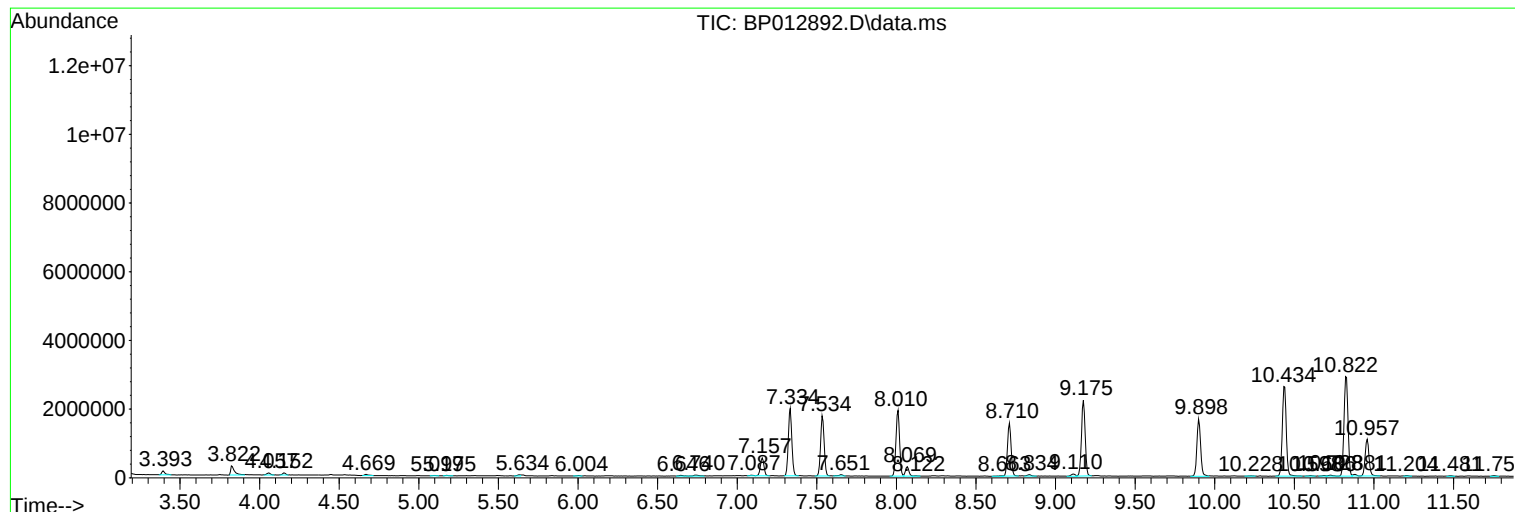
Sum of corrected areas: 148681139

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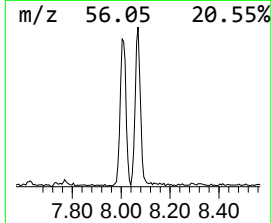
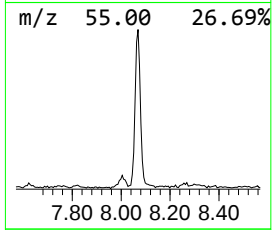
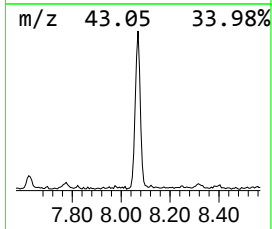
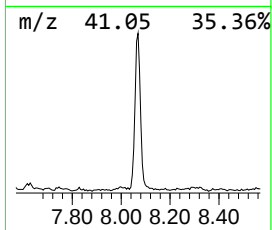
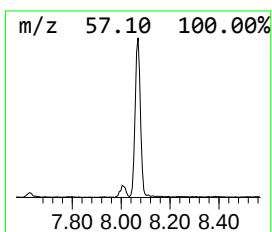
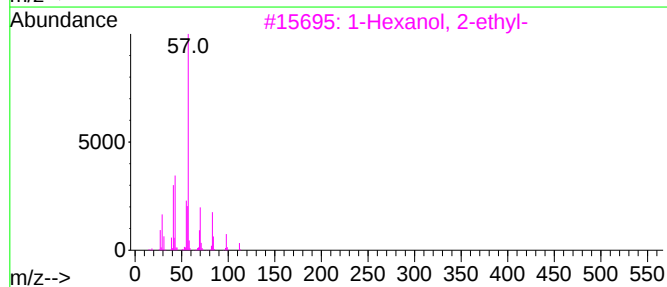
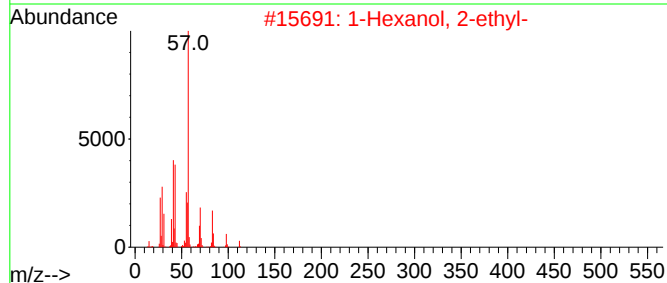
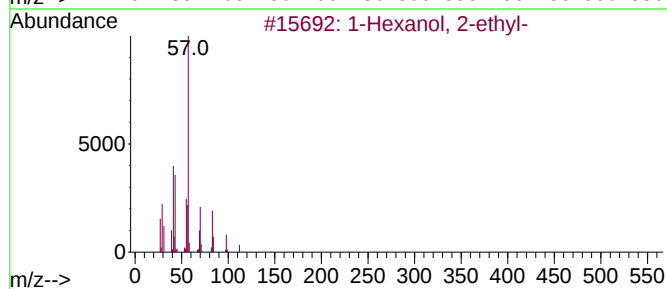
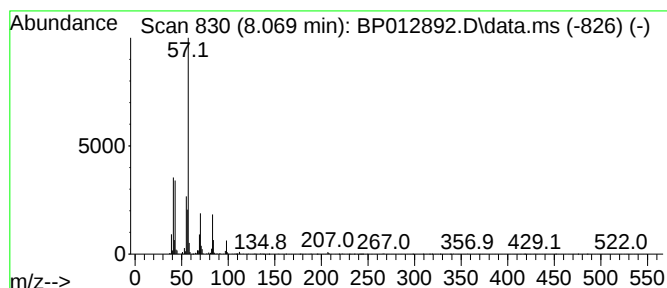
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
Quant Title : SVOA CALIBRATION

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	410899	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
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/u1	410899	1	8.010	3063480	20.0