

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014602.D
 Acq On : 06 Apr 2023 23:29
 Operator : CG/JU
 Sample : 02225-04
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AC9

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014602.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.363	27	30	35	rVV	36919	44684	0.97%	0.101%
2	3.810	103	106	122	rBV	107262	226691	4.93%	0.511%
3	4.016	137	141	148	rVB2	19558	30862	0.67%	0.070%
4	4.116	155	158	165	rVB	11757	18804	0.41%	0.042%
5	4.228	172	177	182	rBV5	10767	17154	0.37%	0.039%
6	4.328	190	194	196	rBV5	3420	4996	0.11%	0.011%
7	5.052	312	317	322	rBV	11724	18259	0.40%	0.041%
8	6.363	536	540	545	rBV5	9400	18997	0.41%	0.043%
9	7.157	666	675	686	rBV	108993	245631	5.34%	0.553%
10	7.310	695	701	716	rBV	581619	947058	20.60%	2.133%
11	7.410	716	718	723	rVB6	3808	4964	0.11%	0.011%
12	7.510	728	735	752	rBV	527776	956755	20.81%	2.155%
13	7.981	808	815	824	rBV	663032	1072313	23.33%	2.416%
14	8.704	931	938	951	rBV	397100	733238	15.95%	1.652%
15	8.992	985	987	993	rVB7	3868	4773	0.10%	0.011%
16	9.157	1008	1015	1032	rBV	725732	1292502	28.12%	2.912%
17	9.287	1032	1037	1050	rVV2	45891	124291	2.70%	0.280%
18	9.522	1075	1077	1086	rVB10	6785	12311	0.27%	0.028%
19	9.887	1131	1139	1156	rBV	612376	1114777	24.25%	2.511%
20	10.428	1222	1231	1251	rBV	703642	1428133	31.07%	3.217%
21	10.798	1286	1294	1307	rBV	948777	1658048	36.07%	3.735%
22	10.951	1313	1320	1339	rBV	487751	1093857	23.79%	2.464%
23	12.022	1498	1502	1507	rVB8	3908	5762	0.13%	0.013%
24	12.404	1560	1567	1576	rBV5	11713	32013	0.70%	0.072%
25	12.539	1585	1590	1599	rVB5	6077	15217	0.33%	0.034%
26	13.445	1735	1744	1748	rBV5	4255	10730	0.23%	0.024%
27	14.039	1836	1845	1856	rBV	1775676	2637412	57.37%	5.941%
28	14.328	1887	1894	1912	rBV	2023356	2976630	64.75%	6.706%
29	14.504	1922	1924	1929	rVB6	3887	5694	0.12%	0.013%
30	14.633	1939	1946	1961	rBV	1580013	2344796	51.01%	5.282%
31	14.922	1988	1995	2004	rBV8	20981	81602	1.78%	0.184%
32	15.463	2083	2087	2090	rBV6	4397	6797	0.15%	0.015%
33	15.627	2108	2115	2130	rBV	2557278	3794421	82.54%	8.548%
34	15.763	2132	2138	2153	rBV	703533	1215612	26.44%	2.738%
35	15.998	2171	2178	2190	rVB	31083	59118	1.29%	0.133%
36	16.333	2230	2235	2245	rVB	4492	13771	0.30%	0.031%

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Peak Location: TOP

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37	16.422	2245	2250	2254	rBV8	3741	7667	0.17%	0.017%
38	17.074	2357	2361	2364	rVB6	5696	8201	0.18%	0.018%
39	17.392	2408	2415	2426	rBV	1850476	2796522	60.83%	6.300%
40	17.492	2426	2432	2446	rVV	2872567	4238058	92.19%	9.547%
41	17.598	2448	2450	2457	rVB6	20051	31754	0.69%	0.072%
42	17.751	2474	2476	2484	rVB6	6032	9662	0.21%	0.022%
43	18.274	2560	2565	2571	rBV6	32898	65365	1.42%	0.147%
44	19.304	2736	2740	2745	rBV3	40477	58240	1.27%	0.131%
45	19.374	2748	2752	2761	rVB2	33383	67364	1.47%	0.152%
46	19.498	2770	2773	2779	rBV8	20116	38197	0.83%	0.086%
47	19.727	2805	2812	2829	rBV	3205567	4597118	100.00%	10.356%
48	21.480	3105	3110	3124	rBV	1496611	2401346	52.24%	5.410%
49	23.821	3500	3508	3520	rBV2	1692031	3647547	79.34%	8.217%
50	23.980	3529	3535	3546	rVB2	998741	2155056	46.88%	4.855%

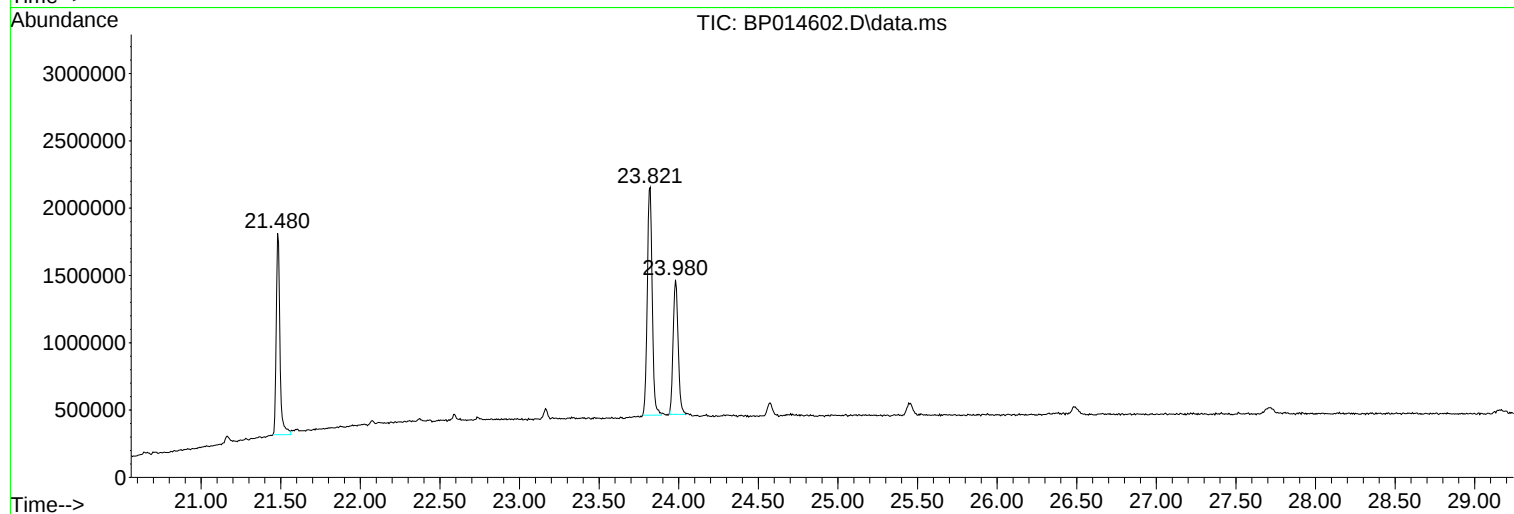
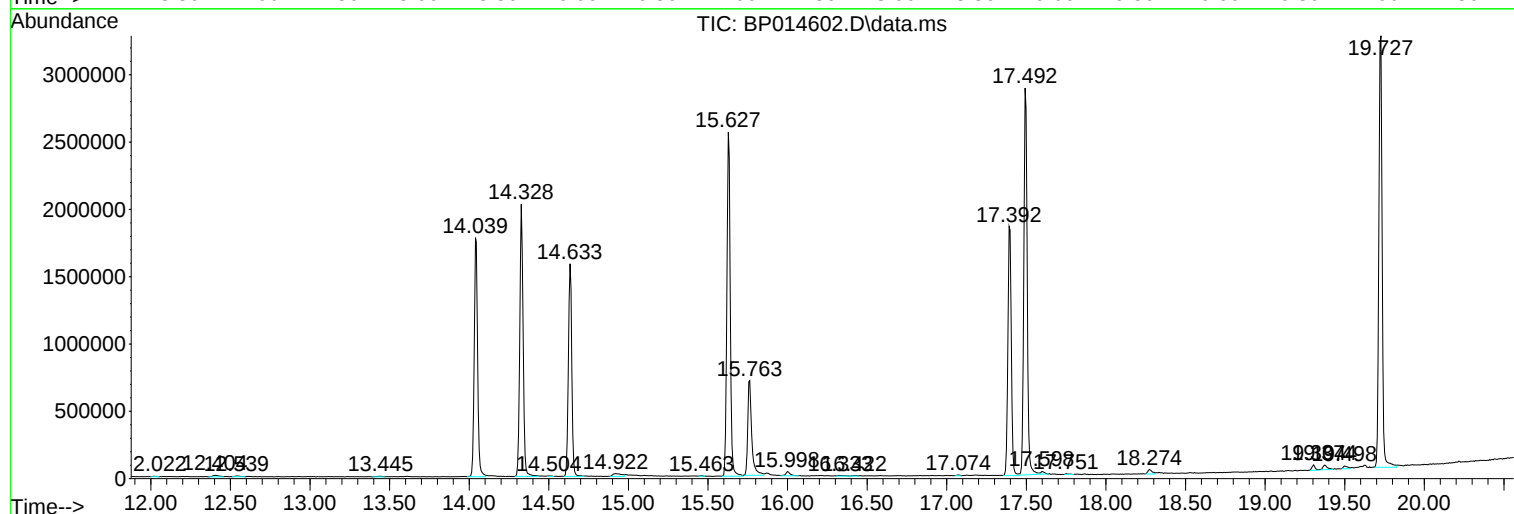
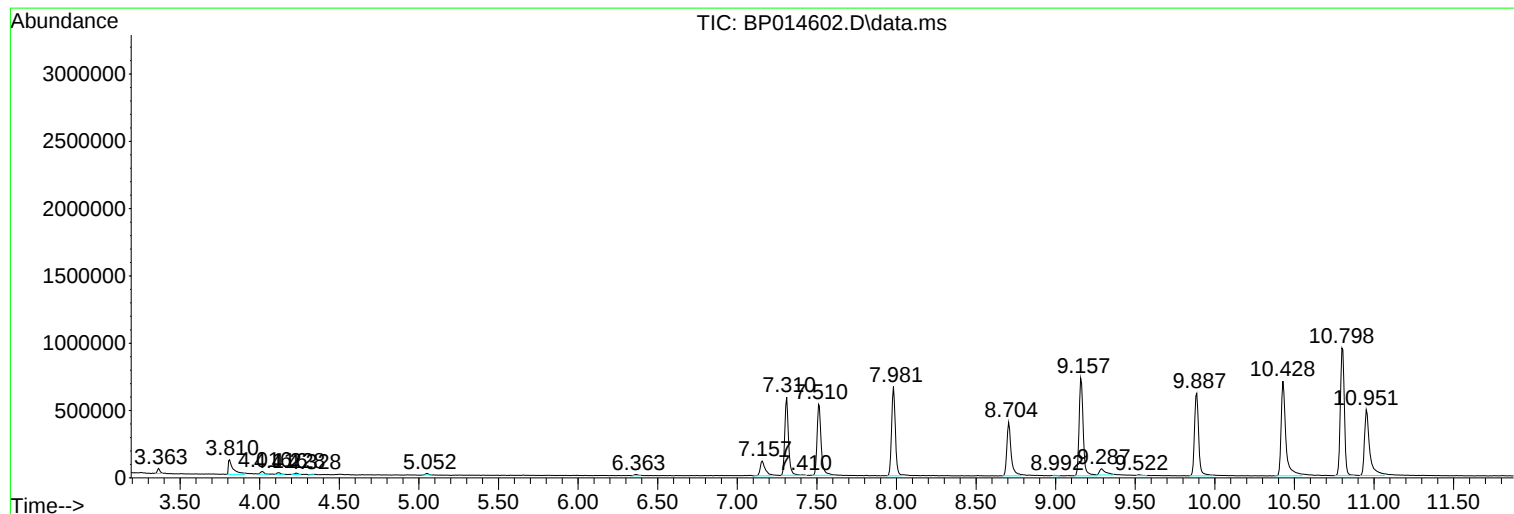
Sum of corrected areas: 44390770

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



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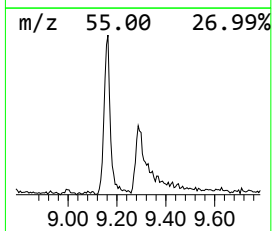
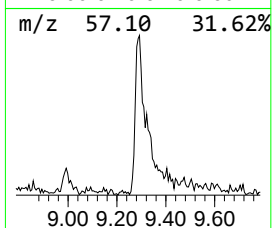
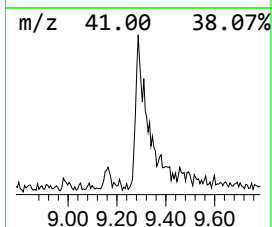
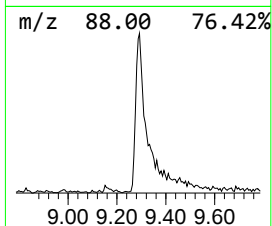
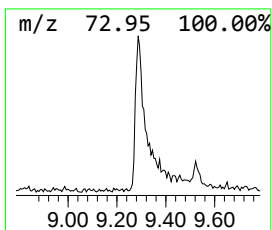
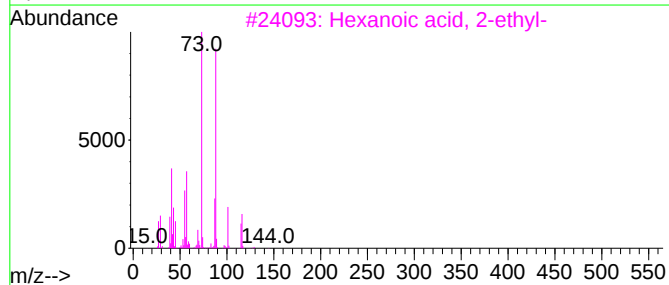
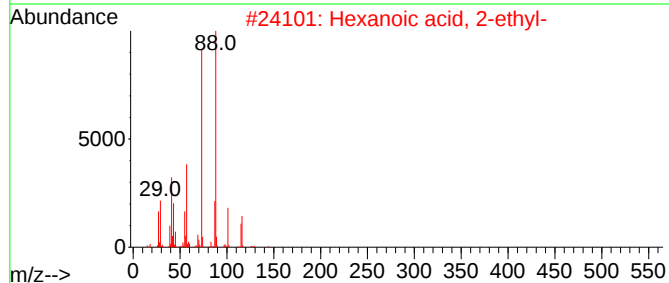
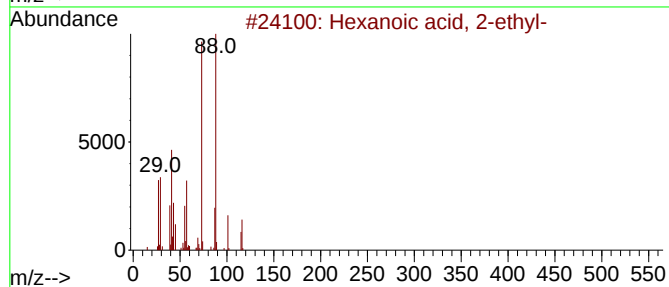
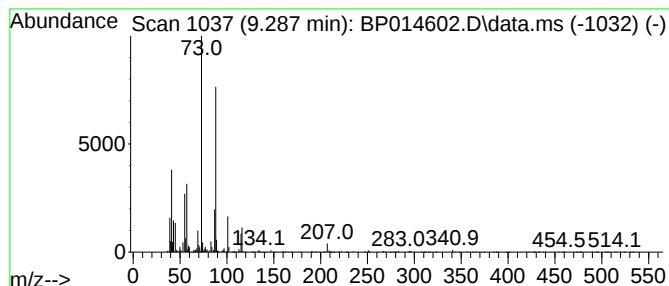
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Hexanoic acid, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.287	2.32 ng/ul	124291	1,4-Dichlorobenzene-d4	7.981

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72



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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
Hexanoic acid, ...	9.287	2.3	ng/ul	124291	1	7.981	1072310	20.0