

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100224\
 Data File : BM047792.D
 Acq On : 03 Oct 2024 01:49
 Operator : RC/JU
 Sample : PB163608BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK608

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BM092724.MA.M

Title : SVOA CALIBRATION

Signal : TIC: BM047792.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	41	45	57	rVB	85988	117652	1.91%	0.167%
2	3.664	112	115	133	rBV	1072827	1593672	25.81%	2.268%
3	3.999	168	172	177	rVB	26413	34223	0.55%	0.049%
4	4.458	245	250	259	rVB	32858	73430	1.19%	0.104%
5	5.505	426	428	432	rBV4	4618	7094	0.11%	0.010%
6	6.728	633	636	641	rVB4	8269	12892	0.21%	0.018%
7	6.969	670	677	692	rBV	1165891	1829355	29.63%	2.603%
8	7.134	699	705	714	rBV	936290	1434372	23.23%	2.041%
9	7.340	732	740	749	rBV	1125823	1797605	29.12%	2.558%
10	7.805	812	819	827	rBV	1106101	1730776	28.03%	2.463%
11	7.875	827	831	836	rVB2	19818	31564	0.51%	0.045%
12	8.040	855	859	861	rBV4	5781	9504	0.15%	0.014%
13	8.510	932	939	955	rBV	1321140	2147062	34.78%	3.055%
14	8.728	971	976	981	rVB4	11335	17420	0.28%	0.025%
15	8.957	1008	1015	1025	rBV	1067022	1733036	28.07%	2.466%
16	9.399	1086	1090	1097	rVB5	17460	31154	0.50%	0.044%
17	9.681	1130	1138	1148	rBV	830663	1358688	22.01%	1.933%
18	10.216	1222	1229	1239	rBV	1271375	2124473	34.41%	3.023%
19	10.328	1247	1248	1254	rVB6	5696	6632	0.11%	0.009%
20	10.598	1285	1294	1302	rBV	1490230	2503380	40.55%	3.562%
21	10.728	1308	1316	1324	rBV	1326230	2205571	35.72%	3.138%
22	11.381	1424	1427	1428	rBV3	5446	6198	0.10%	0.009%
23	11.398	1428	1430	1434	rVB3	5493	6809	0.11%	0.010%
24	11.816	1496	1501	1505	rBV6	7545	12079	0.20%	0.017%
25	11.892	1511	1514	1520	rVB7	5076	7723	0.13%	0.011%
26	12.187	1558	1564	1568	rBV2	20153	34796	0.56%	0.050%
27	12.816	1667	1671	1676	rBV6	7412	14596	0.24%	0.021%
28	13.745	1826	1829	1833	rVB5	6063	7019	0.11%	0.010%
29	13.845	1839	1846	1857	rBV	2324810	3183513	51.56%	4.530%
30	14.134	1883	1895	1903	rBV	2582093	3842183	62.23%	5.467%
31	14.439	1936	1947	1957	rBV	2335601	3327438	53.90%	4.735%
32	14.628	1972	1979	1991	rBV	1050469	1554874	25.18%	2.213%
33	14.845	2014	2016	2021	rVB6	6072	6910	0.11%	0.010%
34	15.251	2081	2085	2090	rVB7	6948	10793	0.17%	0.015%
35	15.428	2108	2115	2122	rBV	3328671	4701258	76.15%	6.690%
36	15.539	2127	2134	2145	rVV	955856	1281182	20.75%	1.823%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	15.669	2153	2156	2161	rVB3	15774	20255	0.33%	0.029%
38	15.869	2187	2190	2195	rVB5	8129	12348	0.20%	0.018%
39	16.116	2227	2232	2233	rVV4	6974	9589	0.16%	0.014%
40	16.133	2233	2235	2240	rVV6	5255	7378	0.12%	0.010%
41	16.204	2244	2247	2253	rVV7	8997	17098	0.28%	0.024%
42	16.504	2293	2298	2304	rBV7	4135	9175	0.15%	0.013%
43	16.827	2348	2353	2356	rBV6	20763	29506	0.48%	0.042%
44	16.969	2372	2377	2379	rVV5	6345	8117	0.13%	0.012%
45	17.051	2387	2391	2397	rBV7	7204	14330	0.23%	0.020%
46	17.175	2405	2412	2422	rBV	2730231	3877452	62.80%	5.518%
47	17.274	2422	2429	2443	rBV	3672579	5139585	83.25%	7.313%
48	17.557	2474	2477	2478	rBV3	6187	6964	0.11%	0.010%
49	17.757	2506	2511	2513	rVB5	5410	7275	0.12%	0.010%
50	18.069	2560	2564	2571	rBV8	15012	33434	0.54%	0.048%
51	18.133	2573	2575	2581	rBV6	4644	6226	0.10%	0.009%
52	18.269	2595	2598	2601	rBV5	4262	6229	0.10%	0.009%
53	18.698	2667	2671	2674	rBV6	9296	14992	0.24%	0.021%
54	19.127	2740	2744	2749	rBV	46308	65036	1.05%	0.093%
55	19.198	2752	2756	2760	rVB2	36894	48583	0.79%	0.069%
56	19.492	2803	2806	2810	rBV4	20493	23689	0.38%	0.034%
57	19.563	2812	2818	2826	rBV	4507917	5974991	96.78%	8.502%
58	20.557	2983	2987	2994	rVB	522326	567944	9.20%	0.808%
59	21.398	3124	3130	3136	rBV	2823194	4202448	68.07%	5.980%
60	22.157	3256	3259	3267	rVB2	136427	235576	3.82%	0.335%
61	24.198	3597	3606	3619	rBV2	2594055	6173828	100.00%	8.785%
62	24.403	3632	3641	3659	rVB	1792919	4966492	80.44%	7.067%

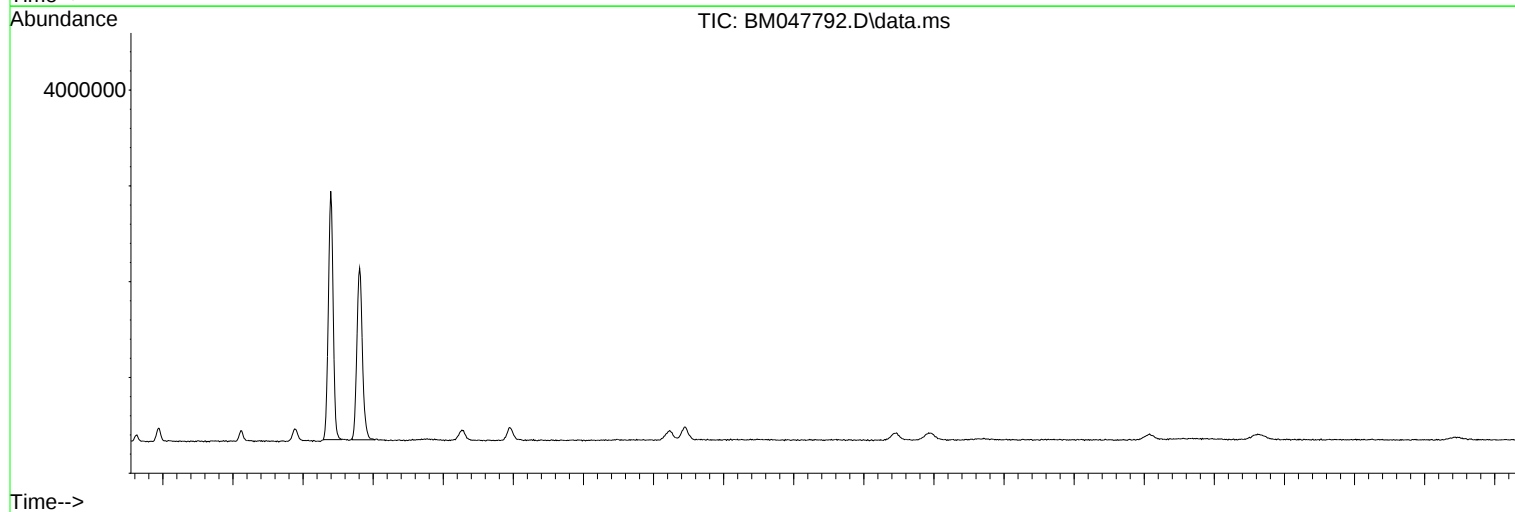
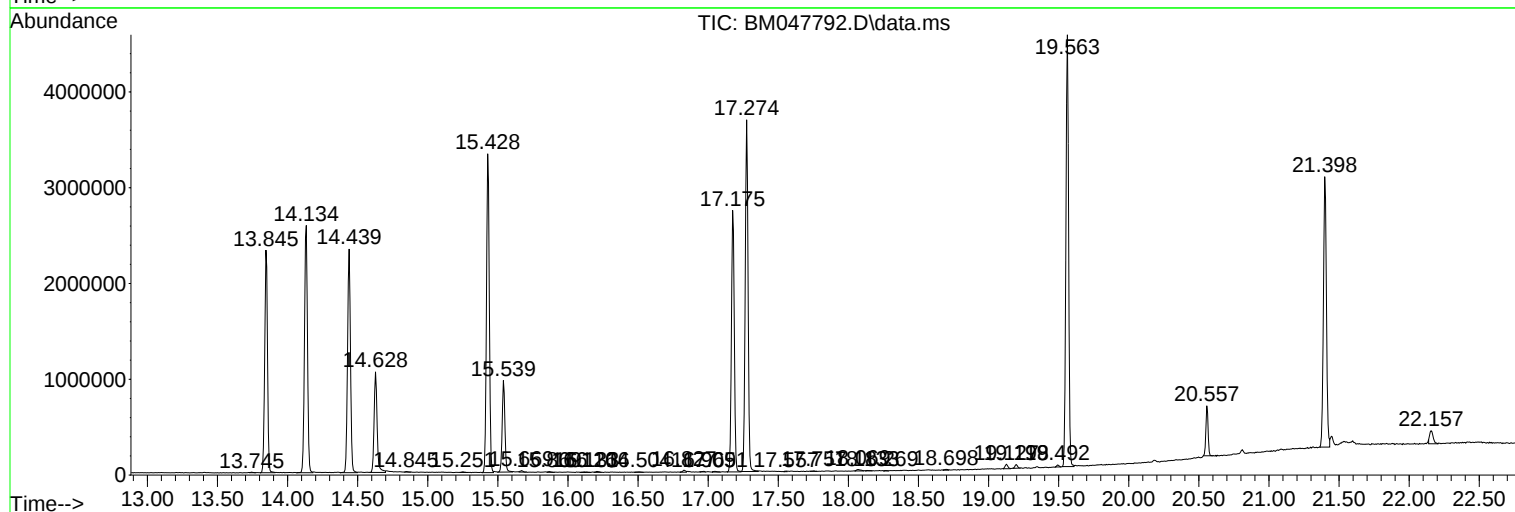
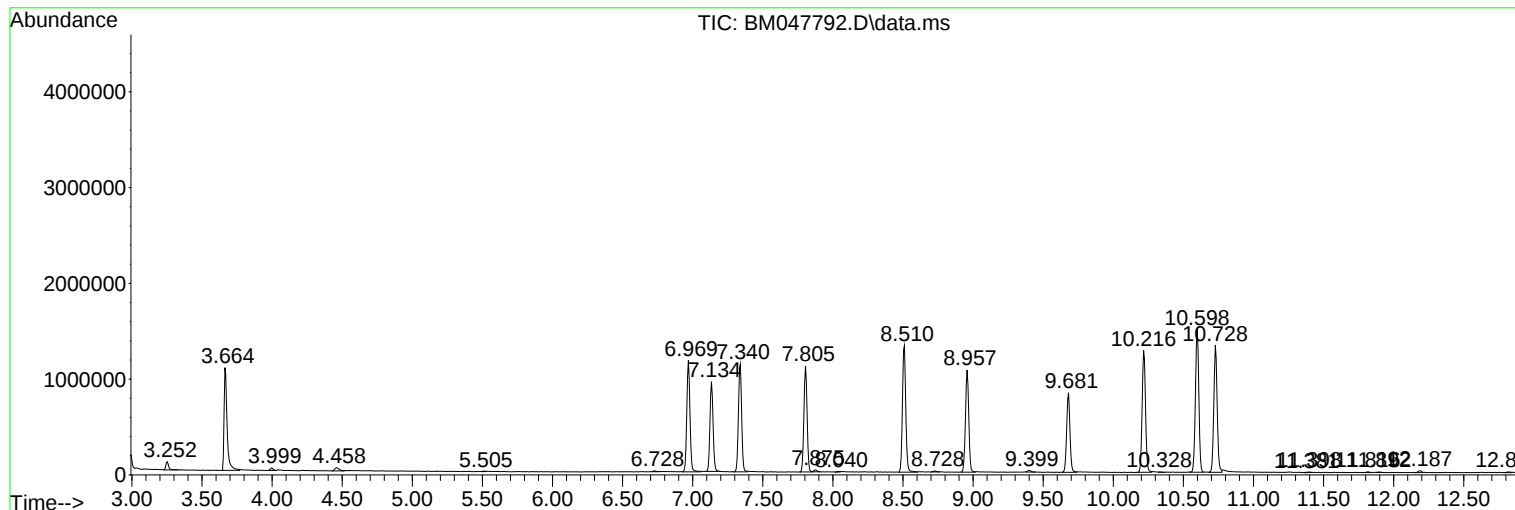
Sum of corrected areas: 70275466

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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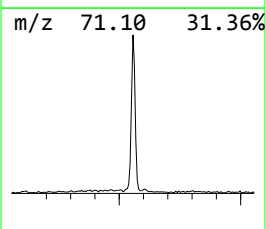
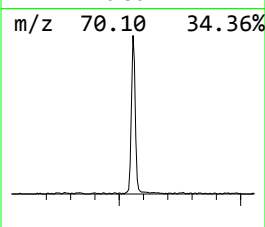
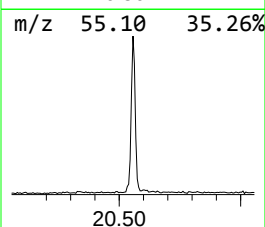
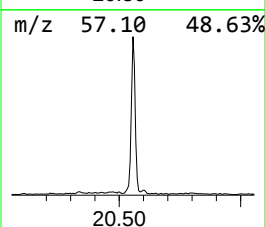
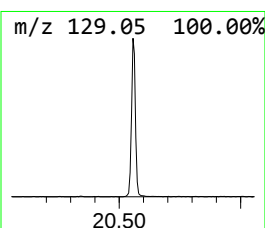
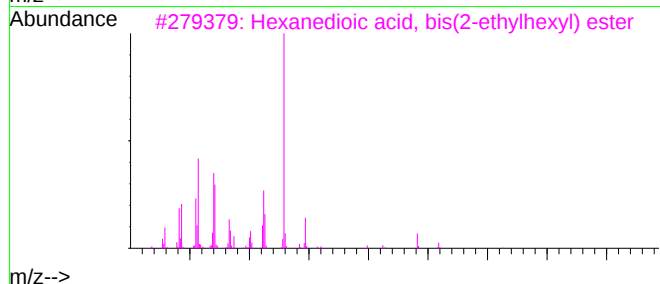
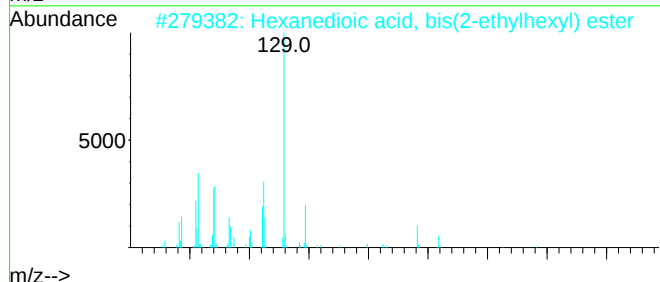
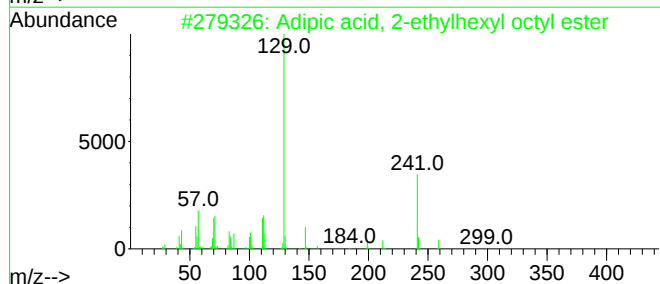
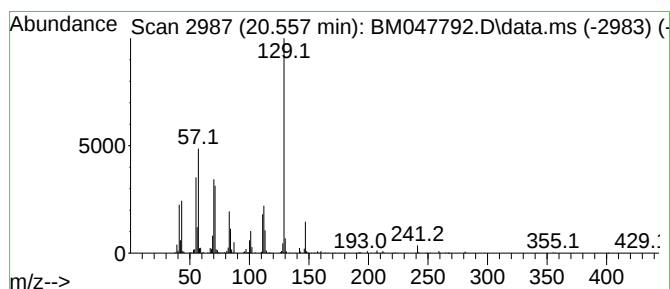
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Adipic acid, 2-ethylhexyl o... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.557	2.70 ng/ul	567944	Chrysene-d12	21.398

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	91
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	86
4	Diisooctyl adipate	370	C22H42O4	001330-86-5	83
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	76



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Adipic acid, 2-...	20.557	2.7	ng/ul	567944	5	21.398	4202450	20.0