

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082021\
 Data File : BN015999.D
 Acq On : 20 Aug 2021 23:17
 Operator : CG/JU
 Sample : M3448-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBKG2

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_N\Methods\SFAM-EPA-BN081621MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN015999.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	109	111	119	rVB2	171493	268220	4.30%	0.452%
2	7.034	666	671	679	rVB	243902	414891	6.65%	0.700%
3	7.205	694	700	711	rVB	730361	1268959	20.35%	2.140%
4	7.405	727	734	742	rBV	824342	1352408	21.69%	2.280%
5	7.875	807	814	821	rVB	876055	1449829	23.25%	2.445%
6	8.581	927	934	944	rVB	632547	1089758	17.48%	1.837%
7	8.810	971	973	978	rVB4	6260	7430	0.12%	0.013%
8	9.034	1003	1011	1019	rBV	1002855	1712818	27.47%	2.888%
9	9.428	1076	1078	1080	rBV3	7200	7019	0.11%	0.012%
10	9.757	1127	1134	1145	rVB	763169	1300722	20.86%	2.193%
11	10.293	1218	1225	1234	rBV	1033677	1809506	29.02%	3.051%
12	10.452	1247	1252	1263	rVV2	37409	71971	1.15%	0.121%
13	10.675	1282	1290	1298	rBV	1156047	2024769	32.48%	3.414%
14	10.810	1305	1313	1327	rVB	1019194	1759894	28.23%	2.967%
15	11.457	1419	1423	1428	rBV6	4769	10242	0.16%	0.017%
16	11.957	1506	1508	1513	rVB5	8168	9983	0.16%	0.017%
17	12.269	1556	1561	1569	rVB3	18743	31756	0.51%	0.054%
18	12.757	1637	1644	1646	rBV6	8996	13478	0.22%	0.023%
19	12.881	1658	1665	1668	rBV7	10733	20252	0.32%	0.034%
20	13.481	1765	1767	1772	rVB3	7007	9108	0.15%	0.015%
21	13.922	1835	1842	1852	rBV	1986306	2949951	47.31%	4.974%
22	14.204	1883	1890	1898	rVB	2339749	3457095	55.45%	5.829%
23	14.410	1922	1925	1929	rBV3	4173	6518	0.10%	0.011%
24	14.510	1936	1942	1952	rBV	1661728	2594001	41.61%	4.374%
25	14.704	1969	1975	1984	rBV	194872	304422	4.88%	0.513%
26	15.251	2064	2068	2073	rBV3	4304	7129	0.11%	0.012%
27	15.316	2075	2079	2081	rBV2	9800	14505	0.23%	0.024%
28	15.504	2101	2111	2119	rBV	2994743	4353091	69.82%	7.340%
29	15.616	2124	2130	2141	rVV	847346	1180045	18.93%	1.990%
30	15.745	2149	2152	2156	rVB3	18260	27357	0.44%	0.046%
31	15.869	2171	2173	2179	rVB5	6579	12280	0.20%	0.021%
32	15.934	2179	2184	2185	rBV4	7603	11252	0.18%	0.019%
33	16.398	2262	2263	2268	rVB5	7437	8873	0.14%	0.015%
34	17.010	2364	2367	2369	rBV3	5052	6271	0.10%	0.011%
35	17.051	2369	2374	2376	rVV5	8641	13359	0.21%	0.023%
36	17.257	2402	2409	2417	rBV	2082422	3046886	48.87%	5.137%

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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_N\Methods\SFAM-EPA-BN081621MA.M
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37	17.357	2419	2426	2433	rBV	3467529	5087857	81.61%	8.578%
38	17.451	2439	2442	2444	rBV4	6768	7449	0.12%	0.013%
39	17.928	2520	2523	2528	rVB6	8896	13018	0.21%	0.022%
40	18.010	2534	2537	2540	rVB4	5156	6702	0.11%	0.011%
41	18.228	2572	2574	2577	rVB	14029	16651	0.27%	0.028%
42	18.457	2609	2613	2615	rBV4	7781	8718	0.14%	0.015%
43	18.569	2629	2632	2634	rBV4	5346	7181	0.12%	0.012%
44	19.022	2705	2709	2711	rBV5	5982	8339	0.13%	0.014%
45	19.086	2717	2720	2722	rBV4	7030	9340	0.15%	0.016%
46	19.216	2738	2742	2748	rBV5	34910	52803	0.85%	0.089%
47	19.280	2749	2753	2763	rVV3	36957	64564	1.04%	0.109%
48	19.451	2777	2782	2785	rBV5	7465	12320	0.20%	0.021%
49	19.480	2785	2787	2791	rVB5	5711	7104	0.11%	0.012%
50	19.510	2791	2792	2794	rBV2	8281	6822	0.11%	0.012%
51	19.533	2794	2796	2799	rBV4	6133	7292	0.12%	0.012%
52	19.651	2810	2816	2822	rBV2	4250896	5963783	95.65%	10.055%
53	19.698	2822	2824	2828	rVB5	22040	22169	0.36%	0.037%
54	19.880	2852	2855	2861	rVB7	13778	23416	0.38%	0.039%
55	20.010	2874	2877	2879	rBV4	9940	11373	0.18%	0.019%
56	20.198	2906	2909	2913	rBV5	24174	29032	0.47%	0.049%
57	20.304	2925	2927	2928	rBV2	9186	6523	0.10%	0.011%
58	20.651	2982	2986	2991	rBV	512385	627273	10.06%	1.058%
59	20.692	2991	2993	2997	rVB2	51409	57702	0.93%	0.097%
60	21.157	3068	3072	3078	rBV4	121038	186700	2.99%	0.315%
61	21.445	3115	3121	3132	rBV	2618983	3507339	56.25%	5.914%
62	21.604	3145	3148	3152	rBV3	58188	65897	1.06%	0.111%
63	22.069	3223	3227	3231	rBV5	62620	82992	1.33%	0.140%
64	22.563	3308	3311	3317	rVB2	89445	121063	1.94%	0.204%
65	23.116	3401	3405	3411	rVB3	112718	168123	2.70%	0.283%
66	23.686	3494	3502	3509	rBV2	2968585	6234731	100.00%	10.512%
67	23.845	3521	3529	3540	rVB2	1642373	3592427	57.62%	6.057%
68	24.463	3630	3634	3643	rVB3	154492	319553	5.13%	0.539%
69	25.310	3773	3778	3786	rVB4	141923	347325	5.57%	0.586%

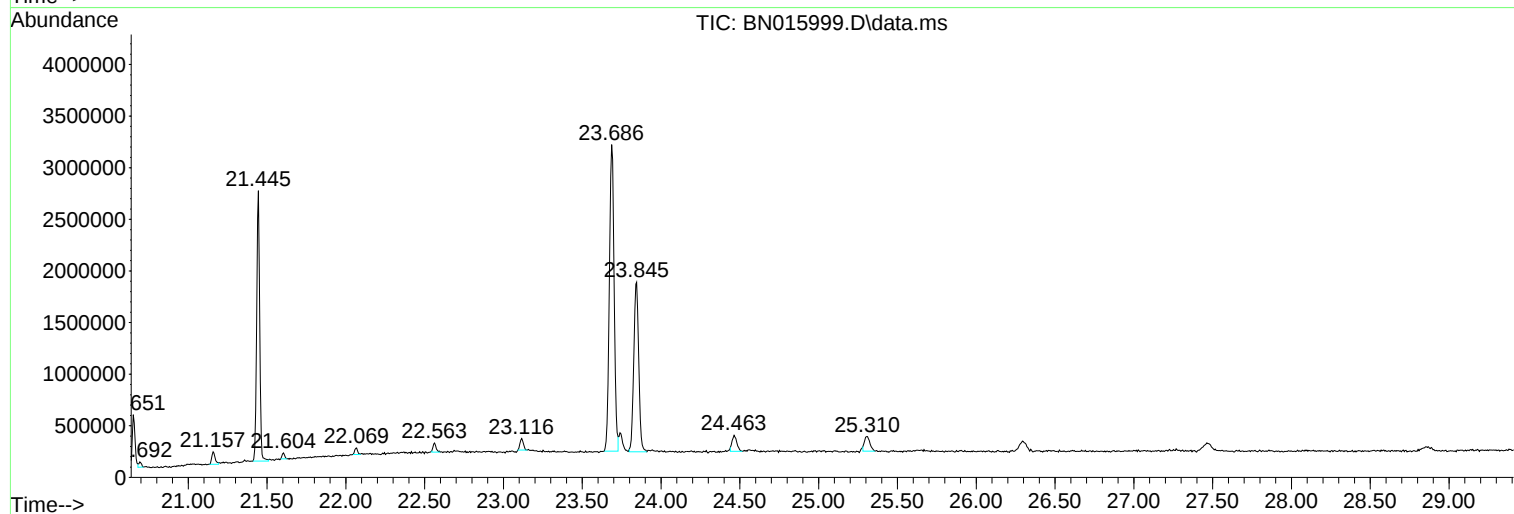
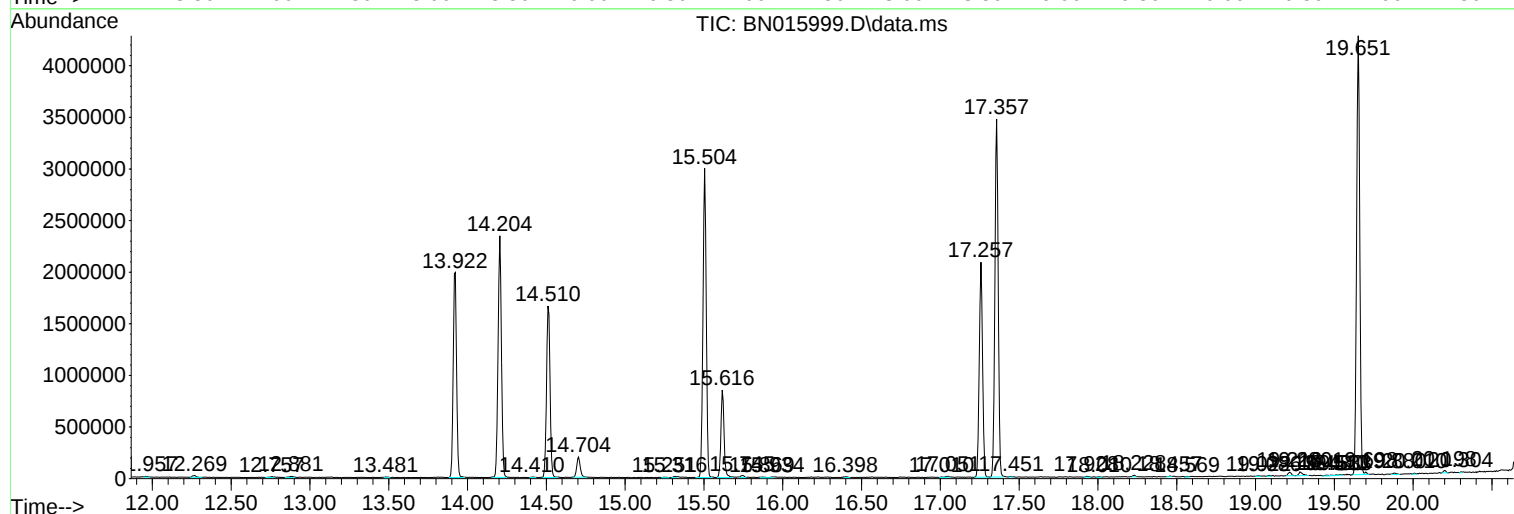
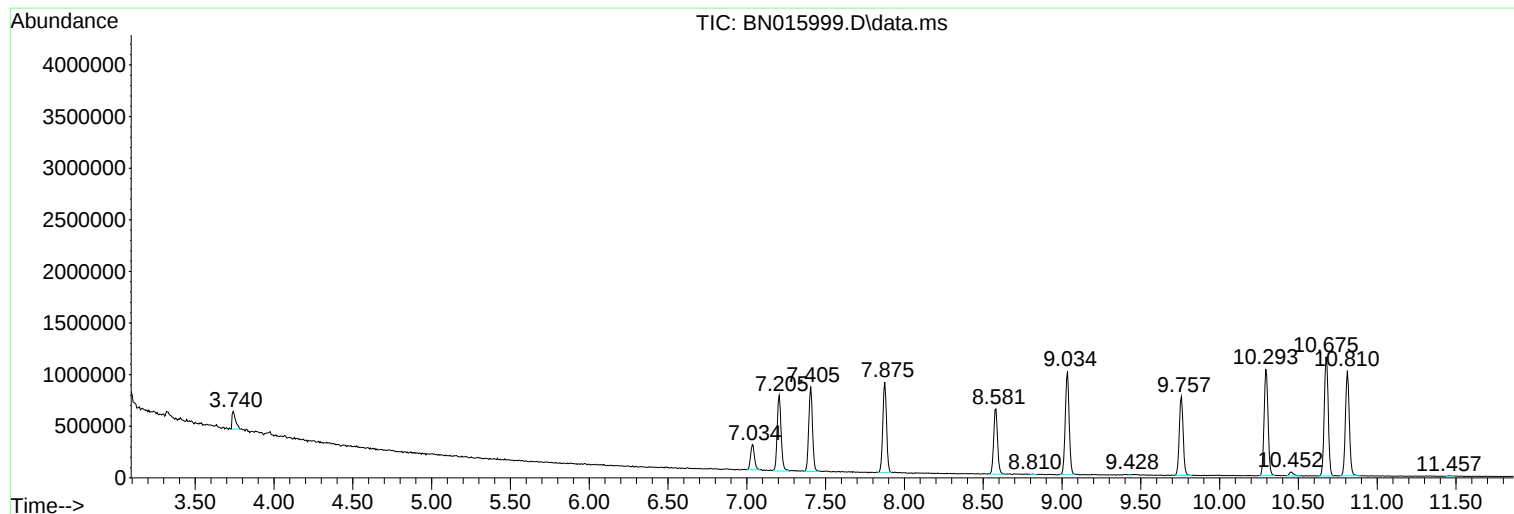
Sum of corrected areas: 59309629

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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



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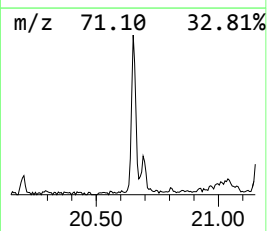
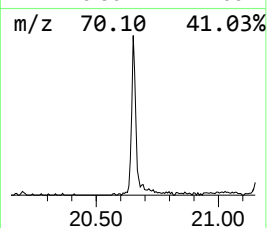
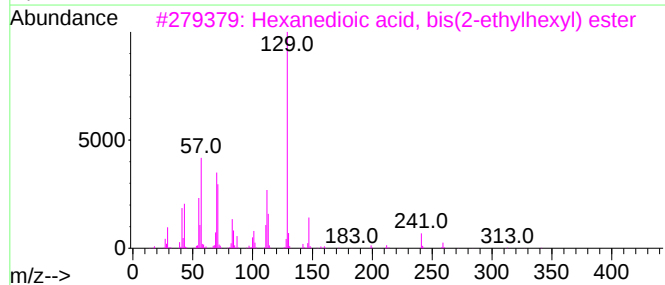
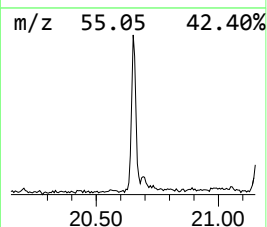
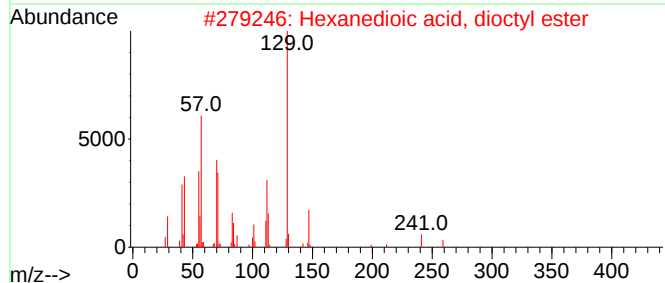
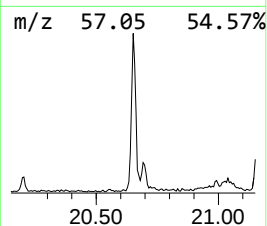
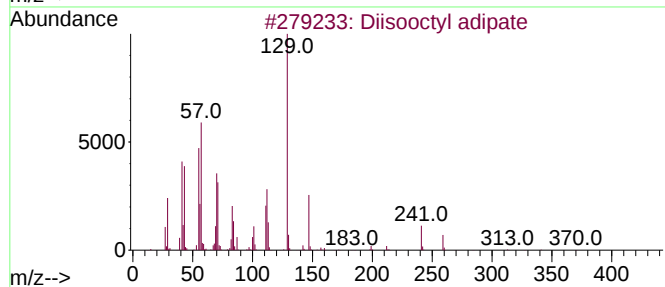
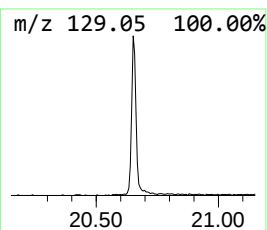
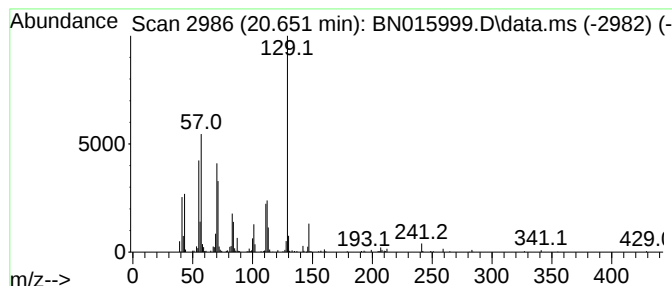
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN081621MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Diisooctyl adipate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.651	3.58 ng/ul	627273	Chrysene-d12	21.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diisooctyl adipate	370	C22H42O4	001330-86-5	91
2			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	90
3			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
4			Adipic acid, 2-ethylhexyl octyl ...	370	C22H42O4	1000324-48-6	90
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Diisooctyl adipate	20.651	3.6	ng/u1	627273	5	21.445	3507340	20.0