

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100820\
 Data File : BN012103.D
 Acq On : 08 Oct 2020 15:30
 Operator : CG/JU
 Sample : PB131925BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK25

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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_N\METHODS\SOM-EPA-BN100220.M
 Title : SVOA CALIBRATION

Signal : TIC: BN012103.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	29	33	42	rVB	83201	122520	1.74%	0.174%
2	3.422	70	74	76	rBV4	6169	7498	0.11%	0.011%
3	3.899	151	155	161	rVB	26790	36979	0.53%	0.052%
4	4.117	189	192	195	rVB4	7214	7559	0.11%	0.011%
5	6.816	645	651	665	rBV	1246801	1943423	27.62%	2.753%
6	6.981	672	679	692	rBV	914548	1468252	20.87%	2.080%
7	7.181	705	713	722	rBV	1281346	2046698	29.09%	2.900%
8	7.640	785	791	799	rBV	909344	1456354	20.70%	2.063%
9	8.346	904	911	922	rBV	1529221	2395010	34.04%	3.393%
10	8.793	979	987	995	rBV	1163432	1893414	26.91%	2.683%
11	9.510	1101	1109	1120	rBV	1007058	1670889	23.75%	2.367%
12	10.040	1191	1199	1211	rBV	1491294	2422379	34.43%	3.432%
13	10.416	1256	1263	1272	rVB	1277099	2075972	29.50%	2.941%
14	10.557	1280	1287	1298	rBV	1419308	2413240	34.30%	3.419%
15	10.657	1302	1304	1308	rVB4	5961	7486	0.11%	0.011%
16	12.010	1530	1534	1541	rBV2	16828	33776	0.48%	0.048%
17	13.257	1741	1746	1749	rBV4	6583	8923	0.13%	0.013%
18	13.698	1815	1821	1833	rBV	2717483	3684061	52.36%	5.219%
19	13.975	1860	1868	1880	rBV	2999701	4302487	61.15%	6.096%
20	14.287	1914	1921	1929	rBV	1924397	2791614	39.68%	3.955%
21	14.487	1949	1955	1975	rBV	1207411	1837862	26.12%	2.604%
22	15.098	2056	2059	2066	rVV6	5479	10757	0.15%	0.015%
23	15.281	2084	2090	2105	rBV	3711435	5202067	73.93%	7.370%
24	15.404	2105	2111	2125	rVV	1380132	1916982	27.25%	2.716%
25	15.522	2126	2131	2136	rVB2	47329	67594	0.96%	0.096%
26	16.069	2220	2224	2227	rVV5	15248	17686	0.25%	0.025%
27	16.822	2349	2352	2354	rBV4	11551	13455	0.19%	0.019%
28	16.957	2372	2375	2381	rBV6	7391	13100	0.19%	0.019%
29	17.034	2381	2388	2395	rBV	2303104	3236200	45.99%	4.585%
30	17.134	2398	2405	2423	rVB	3964946	5623938	79.93%	7.968%
31	17.663	2492	2495	2500	rBV7	6746	8138	0.12%	0.012%
32	17.933	2538	2541	2547	rBV8	25373	35901	0.51%	0.051%
33	18.010	2550	2554	2558	rVB5	21745	31214	0.44%	0.044%
34	18.469	2629	2632	2636	rBV6	19852	25151	0.36%	0.036%
35	18.857	2696	2698	2700	rBV3	7051	7115	0.10%	0.010%
36	18.998	2718	2722	2725	rBV4	22257	29514	0.42%	0.042%

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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

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37	19.069	2731	2734	2737	rVB2	35392	41406	0.59%	0.059%
38	19.280	2768	2770	2772	rBV3	9786	7450	0.11%	0.011%
39	19.322	2773	2777	2780	rBV2	39810	53024	0.75%	0.075%
40	19.433	2790	2796	2805	rBV	4765941	6573472	93.43%	9.313%
41	20.974	3054	3058	3061	rBV	104582	159036	2.26%	0.225%
42	21.022	3062	3066	3071	rBV	386591	497563	7.07%	0.705%
43	21.233	3097	3102	3110	rBV2	2868980	3674582	52.23%	5.206%
44	23.298	3447	3453	3468	rVB	3700506	7036010	100.00%	9.968%
45	23.433	3471	3476	3484	rVB	2046995	3676105	52.25%	5.208%

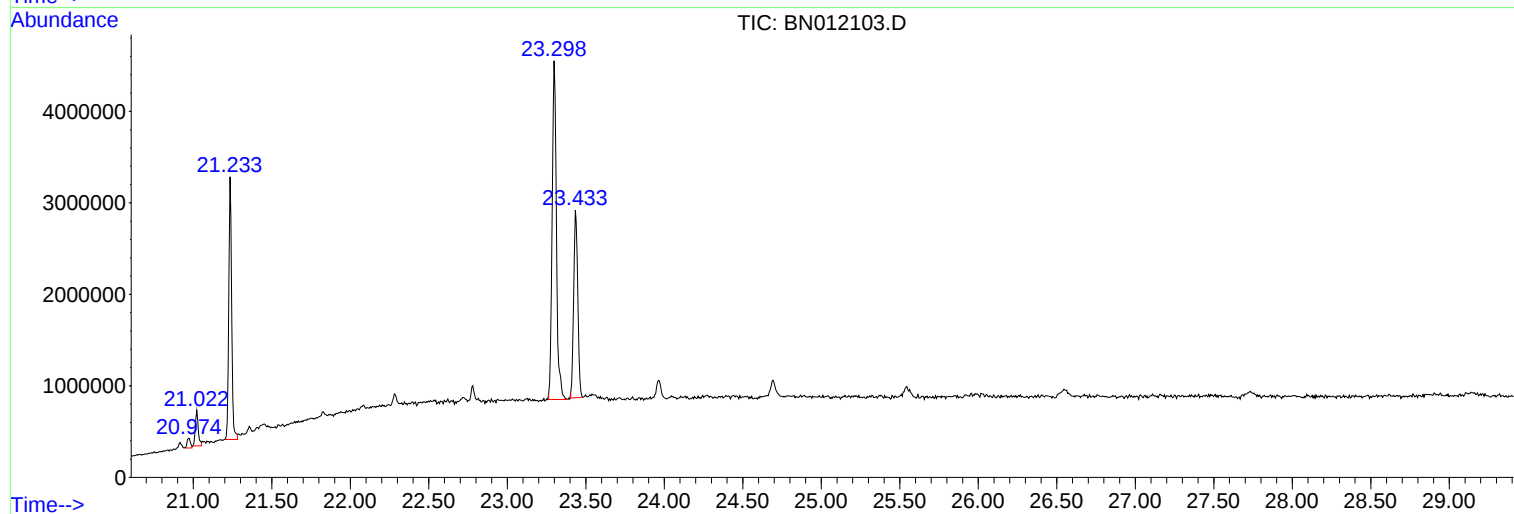
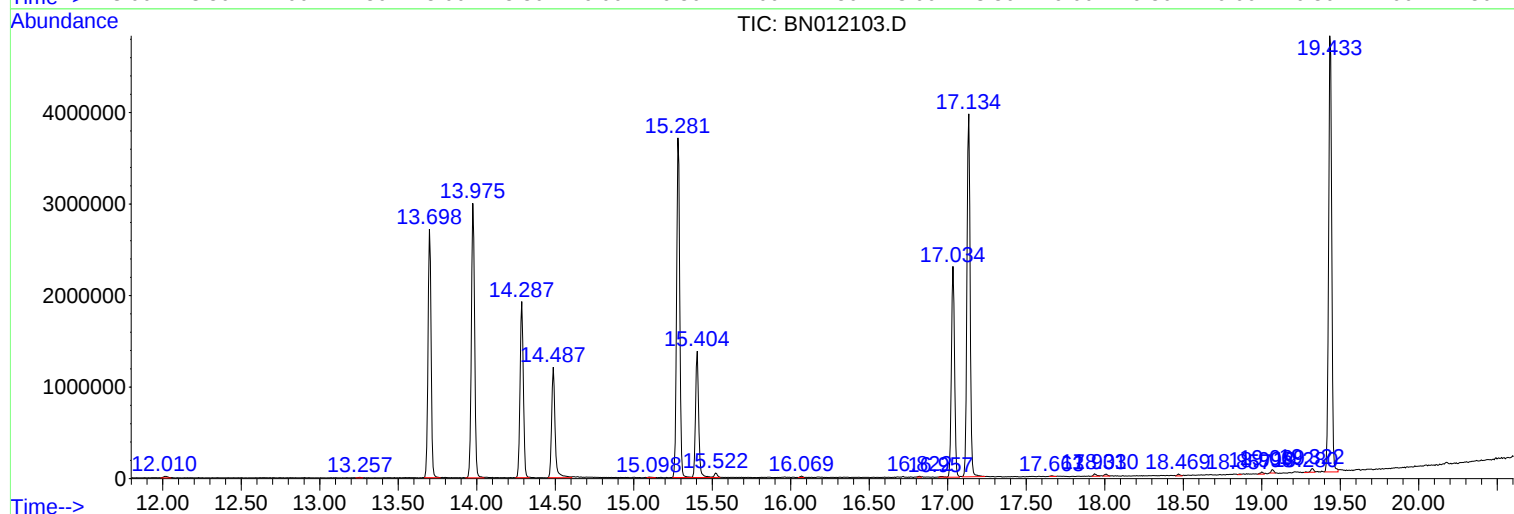
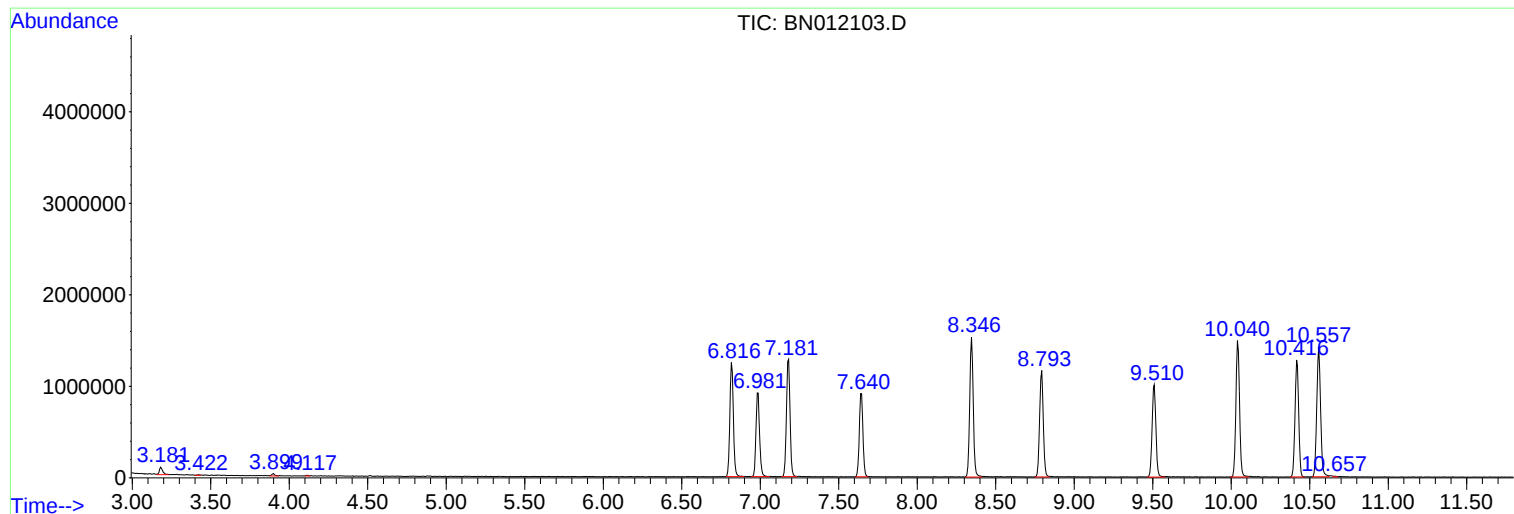
Sum of corrected areas: 70583856

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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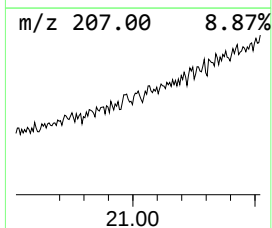
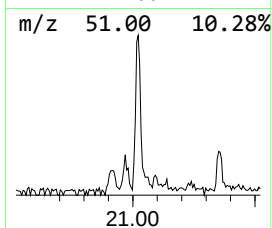
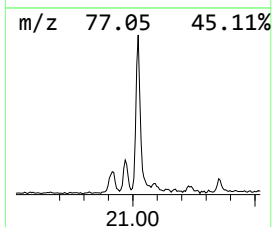
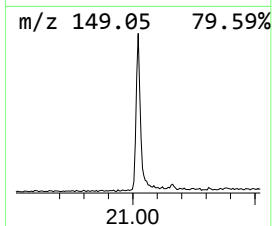
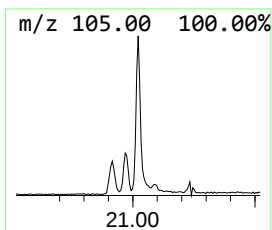
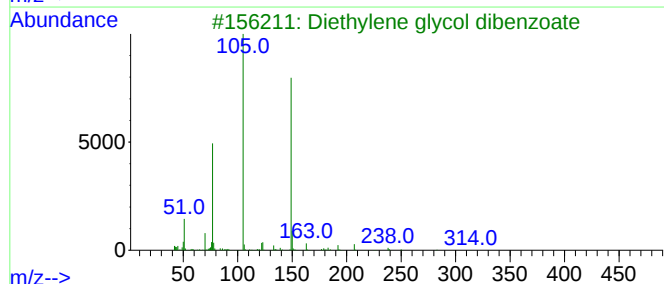
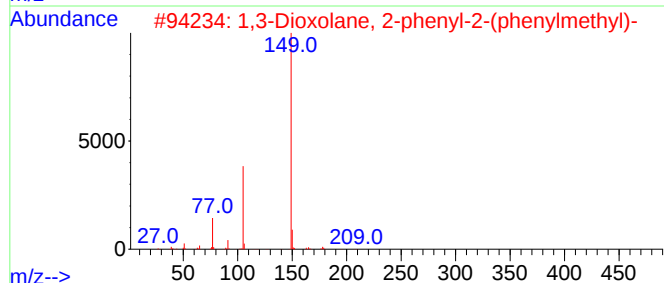
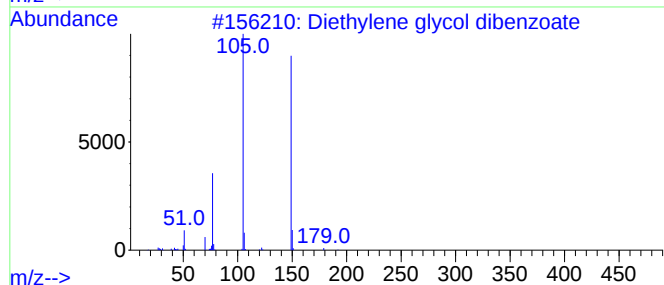
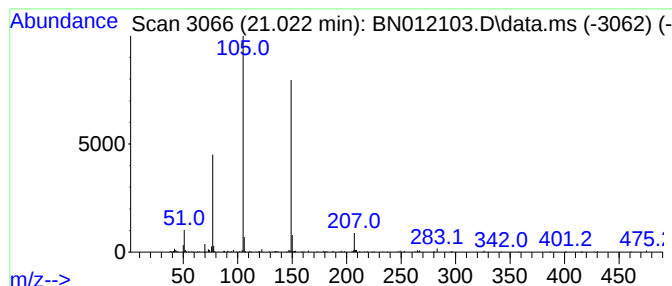
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.020	2.71 ng/ul	497563	Chrysene-d12	21.233

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	86
2			1,3-Dioxolane, 2-phenyl-2-(pheny...	240	C16H16O2	004362-19-0	78
3			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	64
4			2-Bromomethyl-2-phenyl[1,3]dioxo...	242	C10H11BrO2	003418-21-1	64
5			Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	56



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Diethylene glyc...	21.020	2.7	ng/u1	497563	5	21.233	3674580	20.0