

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112520\
 Data File : BN012738.D
 Acq On : 25 Nov 2020 19:06
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL-01
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK233

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_N\METHODS\SFAM-EPA-BN112520.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.617	20	22	27	rVB2	19415	24128	0.54%	0.053%
2	2.699	33	36	41	rVB2	16058	23034	0.52%	0.050%
3	2.958	77	80	87	rVB	53463	62378	1.40%	0.136%
4	3.358	146	148	166	rBV	653050	1049629	23.55%	2.288%
5	3.669	198	201	209	rVB	17889	24704	0.55%	0.054%
6	6.610	693	701	713	rBV	877648	1312166	29.44%	2.860%
7	6.775	723	729	739	rBV	612244	911861	20.46%	1.988%
8	6.957	753	760	772	rBV	910432	1365352	30.63%	2.976%
9	7.422	832	839	850	rBV	512426	799655	17.94%	1.743%
10	8.134	954	960	975	rBV	1005143	1562119	35.04%	3.405%
11	8.575	1028	1035	1051	rBV	797805	1316120	29.53%	2.869%
12	9.287	1149	1156	1168	rBV	742045	1167486	26.19%	2.545%
13	9.816	1239	1246	1257	rBV	1002778	1613934	36.21%	3.518%
14	10.187	1302	1309	1318	rBV	704746	1098713	24.65%	2.395%
15	10.334	1328	1334	1354	rBV	916114	1623007	36.41%	3.538%
16	11.793	1576	1582	1583	rBV3	8809	9040	0.20%	0.020%
17	13.504	1867	1873	1881	rBV	1871502	2459387	55.17%	5.361%
18	13.769	1912	1918	1929	rBV	1937760	2772062	62.19%	6.042%
19	14.081	1965	1971	1981	rBV	1003206	1479163	33.18%	3.224%
20	14.304	2003	2009	2023	rBV	840146	1258821	28.24%	2.744%
21	15.086	2135	2142	2151	rBV	2436557	3335872	74.84%	7.271%
22	15.222	2159	2165	2173	rBV	942410	1291367	28.97%	2.815%
23	15.322	2178	2182	2188	rVB	24372	36888	0.83%	0.080%
24	15.875	2270	2276	2281	rBV7	8037	11531	0.26%	0.025%
25	16.622	2400	2403	2406	rBV3	5361	6272	0.14%	0.014%
26	16.839	2434	2440	2450	rBV	1290335	1738669	39.01%	3.790%
27	16.933	2450	2456	2471	rBV2	2477097	3629865	81.43%	7.912%
28	17.763	2591	2597	2600	rBV6	5045	8237	0.18%	0.018%
29	18.369	2696	2700	2702	rBV3	3916	6983	0.16%	0.015%
30	18.816	2773	2776	2781	rBV6	4240	6817	0.15%	0.015%
31	18.880	2783	2787	2791	rVV3	20527	27313	0.61%	0.060%
32	19.039	2811	2814	2815	rBV2	5338	4968	0.11%	0.011%
33	19.133	2826	2830	2835	rVB	19514	25074	0.56%	0.055%
34	19.245	2843	2849	2862	rVB2	3085388	4194662	94.10%	9.143%

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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

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Title : SVOA CALIBRATION

35	20.751	3099	3105	3109	rBV	114437	199806	4.48%	0.436%
36	20.804	3110	3114	3118	rVV	168817	219768	4.93%	0.479%
37	20.851	3118	3122	3130	rVV	462801	620096	13.91%	1.352%
38	21.057	3152	3157	3164	rBV	1578236	1929468	43.29%	4.206%
39	21.180	3175	3178	3183	rBV4	47408	81669	1.83%	0.178%
40	23.039	3487	3494	3499	rBV	2720219	4457476	100.00%	9.716%
41	23.168	3510	3516	3525	rVB	1209941	2112226	47.39%	4.604%

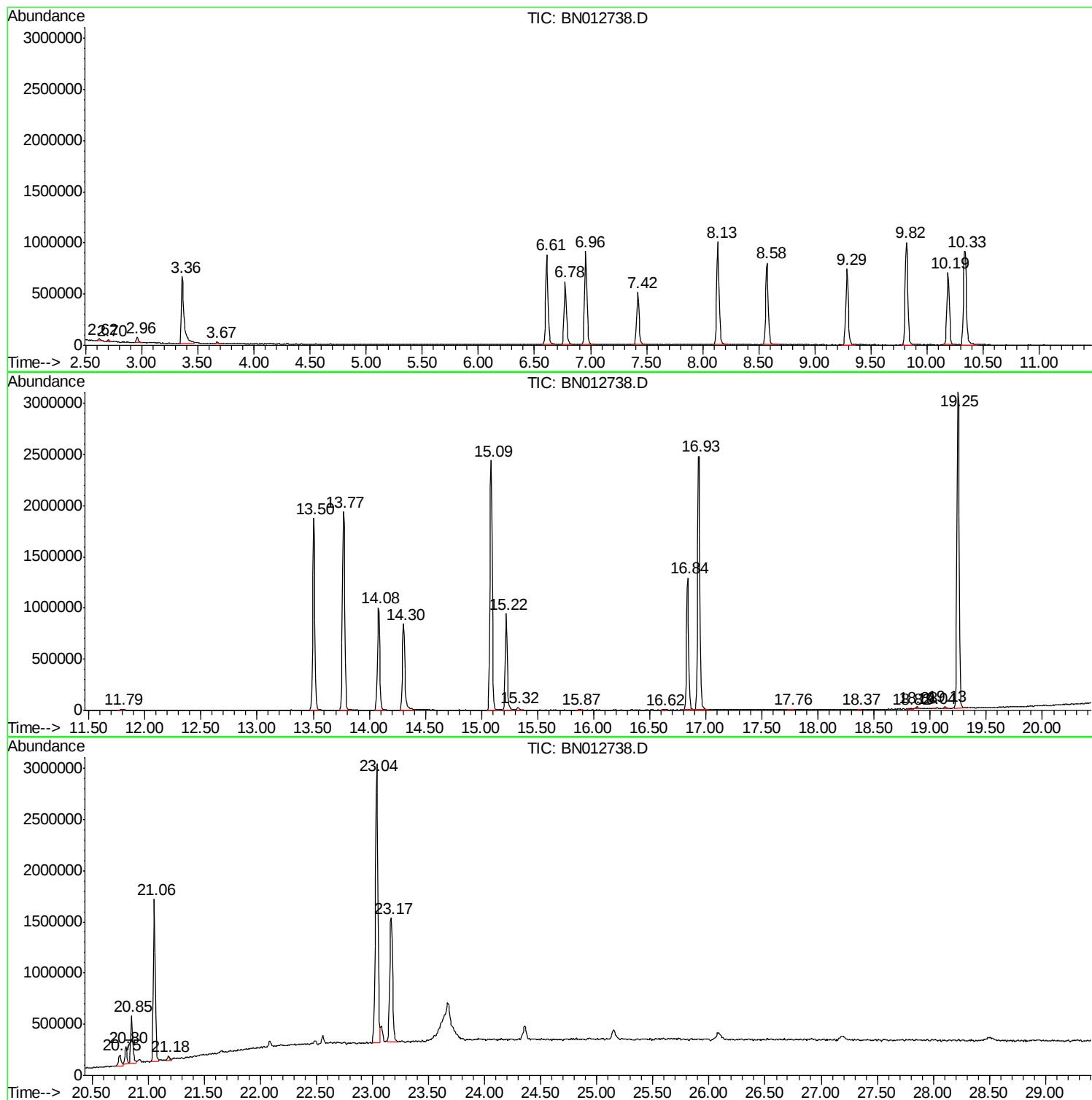
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN112520.M
Quant Title : SVOA CALIBRATION

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



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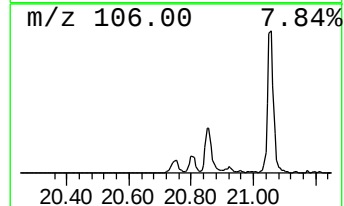
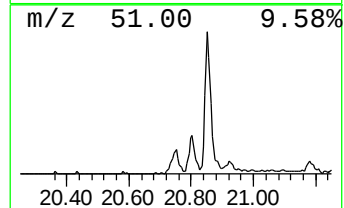
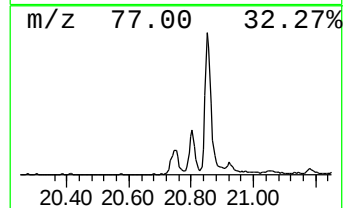
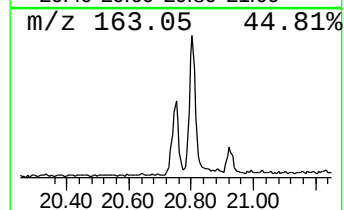
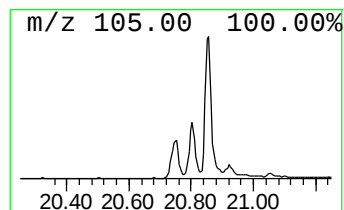
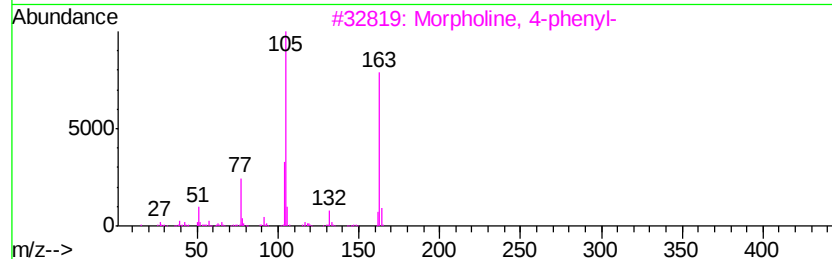
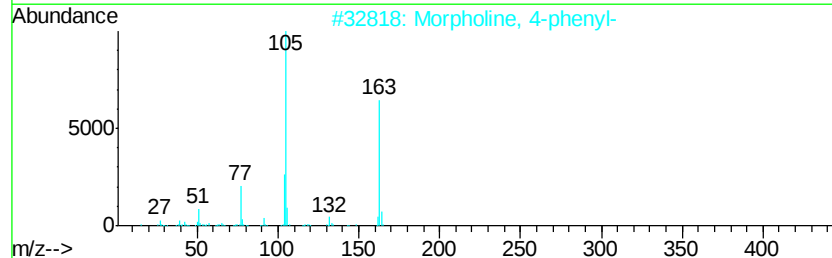
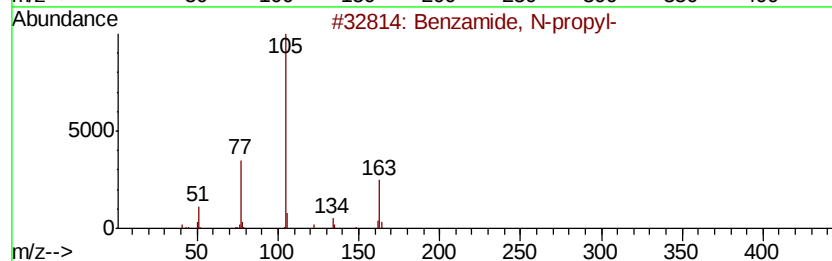
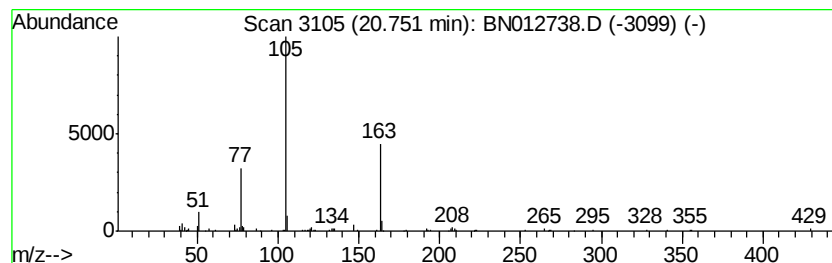
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN112520.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzamide, N-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.75	2.07 ng/ul	199806	Chrysene-d12	21.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, N-propyl-	163	C10H13NO	010546-70-0	78
2		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64
3		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
4		Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	38
5		Glyoxylic acid, phenyl-, 2'-cyan...	251	C15H9NO3	166538-14-3	38



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Misc : MDL-WATER-BL-01
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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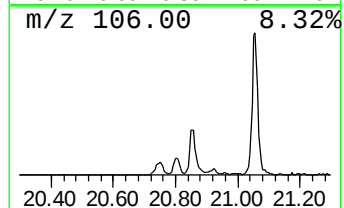
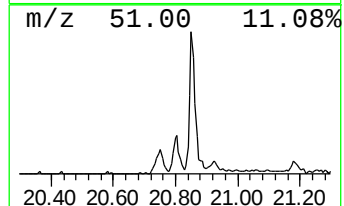
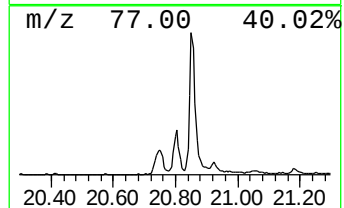
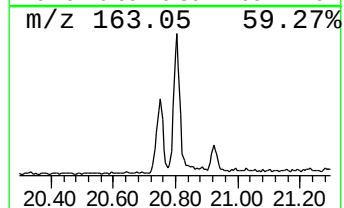
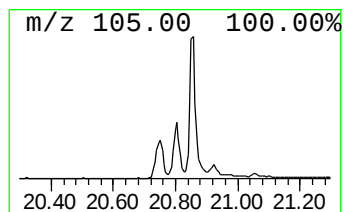
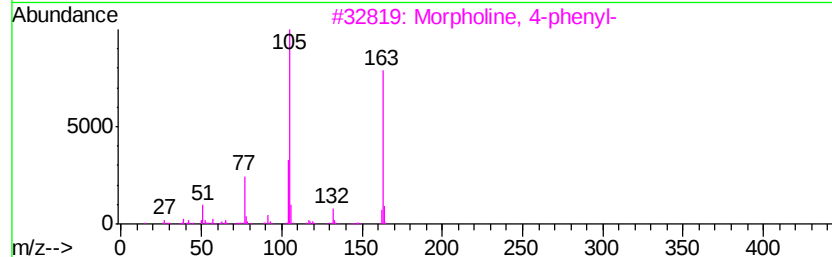
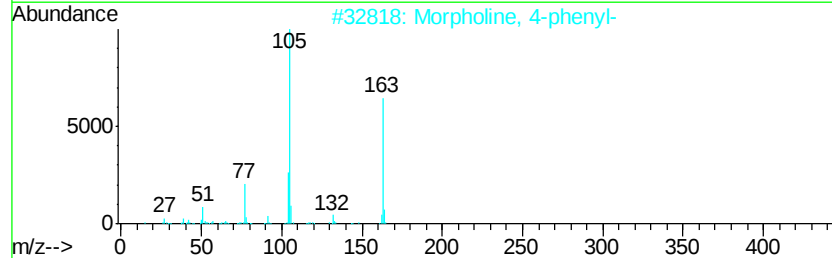
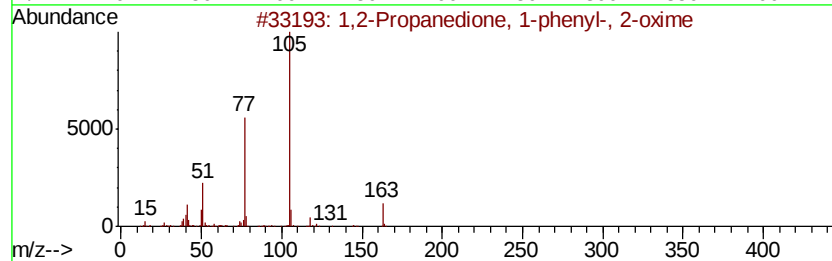
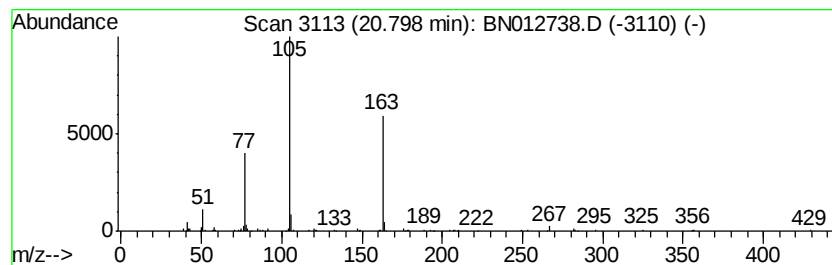
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Peak Number 2 1,2-Propanedione, 1-phenyl-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.80	2.28 ng/ul	219768	Chrysene-d12	21.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	50
2		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	47
3		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	47
4		Benzaldehyde, 4-cyano-, benzoylh...	249	C15H11N3O	195066-12-7	43
5		1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	43



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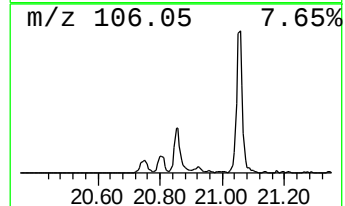
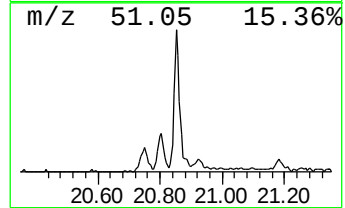
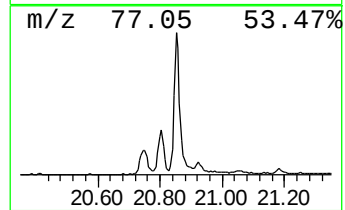
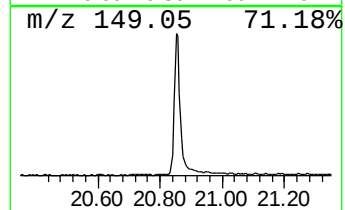
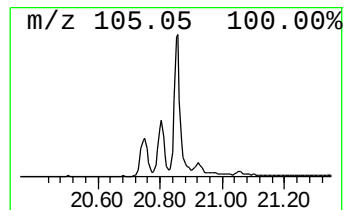
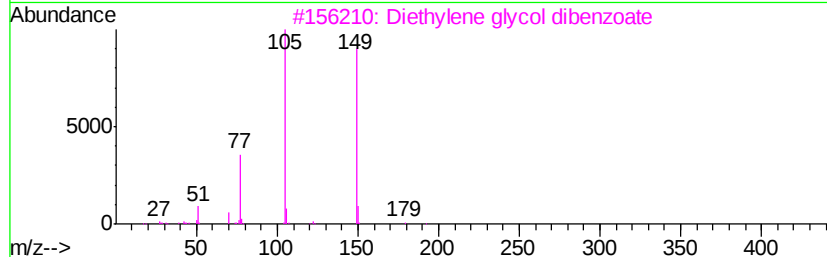
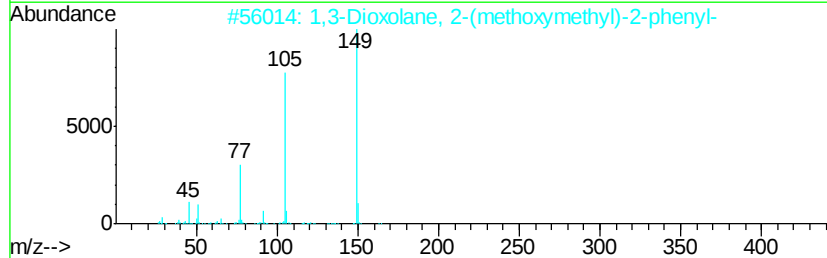
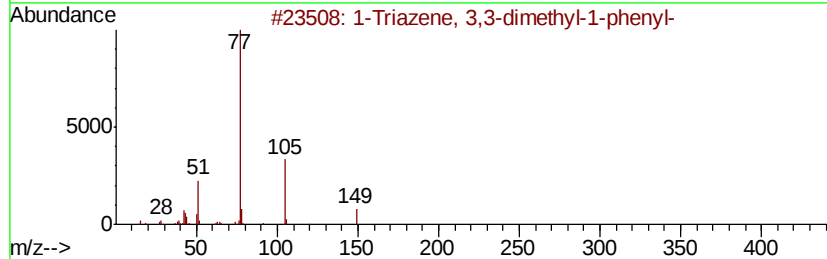
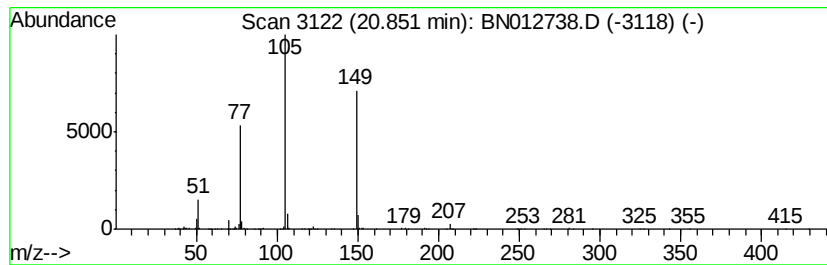
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Triazene, 3,3-dimethyl-1-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.		
20.85	6.43 ng/ul	620096	Chrysene-d12		21.06		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Triazene, 3,3-dimethyl-1-phenyl-	149	C8H11N3	007227-91-0	72		
2	1,3-Dioxolane, 2-(methoxymethyl)-2-phenyl-	194	C11H14O3	1000156-71-2	53		
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	50		
4	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	46		
5	2-Benzoyloxyacetophenone	240	C15H12O3	004010-33-7	43		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzamide, N-propyl-	20.75	2.1	ng/ul	199806	5	21.06	1929470	20.0
1,2-Propanedione,...	20.80	2.3	ng/ul	219768	5	21.06	1929470	20.0
1-Triazene, 3,3-d...	20.85	6.4	ng/ul	620096	5	21.06	1929470	20.0