

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
 Data File : BM026333.D
 Acq On : 10 Jun 2020 17:01
 Operator : JU/CG
 Sample : L2884-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BFXD7

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\SOM-EPA-BM051320MA.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	3.005	17	20	26	rBV	58813	73615	2.19%	0.155%
2	3.070	27	31	35	rVB2	27104	41965	1.25%	0.088%
3	4.922	340	346	356	rVV	284032	387049	11.53%	0.813%
4	5.134	373	382	387	rBV2	14225	31383	0.93%	0.066%
5	5.469	431	439	443	rBV	58687	83773	2.50%	0.176%
6	6.628	625	636	648	rVB2	177225	361443	10.77%	0.759%
7	6.775	654	661	668	rBV	592397	907632	27.03%	1.907%
8	6.963	686	693	701	rBV	668880	1011346	30.12%	2.125%
9	7.093	710	715	720	rBV2	32906	50461	1.50%	0.106%
10	7.422	764	771	779	rBV	694779	1055452	31.44%	2.217%
11	7.505	779	785	787	rVV4	16575	32928	0.98%	0.069%
12	7.534	787	790	800	rVB2	27031	47756	1.42%	0.100%
13	7.769	822	830	838	rVB2	35068	75380	2.25%	0.158%
14	8.146	887	894	902	rVV	479287	751923	22.40%	1.580%
15	8.422	934	941	948	rBV	60571	88436	2.63%	0.186%
16	8.569	959	966	975	rBV	701675	1127811	33.59%	2.369%
17	8.681	975	985	989	rVB4	25991	55253	1.65%	0.116%
18	9.034	1039	1045	1050	rVB3	21341	36718	1.09%	0.077%
19	9.281	1079	1087	1097	rBV	630519	1051346	31.31%	2.209%
20	9.446	1110	1115	1121	rVB2	47523	78102	2.33%	0.164%
21	9.593	1131	1140	1144	rBV2	22288	39027	1.16%	0.082%
22	9.816	1172	1178	1188	rBV	801904	1364058	40.63%	2.866%
23	10.087	1219	1224	1229	rVV	27860	43706	1.30%	0.092%
24	10.181	1229	1240	1246	rVV	885872	1426705	42.49%	2.997%
25	10.228	1246	1248	1253	rVB	33503	45468	1.35%	0.096%
26	10.334	1258	1266	1271	rBV2	26052	60050	1.79%	0.126%
27	10.604	1302	1312	1317	rBV8	16442	39196	1.17%	0.082%
28	10.793	1340	1344	1355	rVB10	14158	33527	1.00%	0.070%
29	11.181	1404	1410	1412	rBV7	20202	36949	1.10%	0.078%
30	11.281	1424	1427	1434	rVV3	19125	38037	1.13%	0.080%
31	11.357	1434	1440	1451	rVV	110808	184172	5.49%	0.387%
32	11.457	1453	1457	1463	rVB6	19256	31229	0.93%	0.066%
33	11.522	1463	1468	1478	rBV4	52489	98391	2.93%	0.207%
34	11.751	1498	1507	1508	rBV8	23723	50970	1.52%	0.107%

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35	11.781	1509	1512	1517	rVV2	46271	68690	2.05%	0.144%
36	11.834	1517	1521	1525	rVB5	24294	34839	1.04%	0.073%
37	11.887	1525	1530	1536	rBV6	17941	31384	0.93%	0.066%
38	12.040	1549	1556	1559	rVB5	18909	39541	1.18%	0.083%
39	12.175	1573	1579	1582	rVV4	26944	56523	1.68%	0.119%
40	12.316	1590	1603	1607	rBV	42700	81992	2.44%	0.172%
41	12.445	1619	1625	1629	rVV7	23790	42015	1.25%	0.088%
42	12.551	1637	1643	1649	rBV2	143194	230727	6.87%	0.485%
43	12.628	1649	1656	1658	rVV2	29715	62749	1.87%	0.132%
44	12.704	1664	1669	1676	rVB7	17873	39445	1.17%	0.083%
45	12.898	1694	1702	1713	rBV4	32671	104355	3.11%	0.219%
46	13.122	1734	1740	1748	rVV	389569	540363	16.09%	1.135%
47	13.222	1751	1757	1760	rBV7	18565	37399	1.11%	0.079%
48	13.257	1760	1763	1768	rVB2	23004	31949	0.95%	0.067%
49	13.498	1797	1804	1809	rBV	1372917	1938733	57.74%	4.073%
50	13.545	1809	1812	1817	rVB	264329	331831	9.88%	0.697%
51	13.757	1837	1848	1856	rBV	1615079	2606249	77.62%	5.475%
52	14.075	1896	1902	1908	rVV	1212606	1746839	52.03%	3.670%
53	14.128	1908	1911	1917	rVB3	46675	68226	2.03%	0.143%
54	14.322	1939	1944	1955	rVB3	101781	218070	6.49%	0.458%
55	14.445	1959	1965	1970	rBV3	56930	117321	3.49%	0.246%
56	14.639	1994	1998	2004	rVV5	35482	66593	1.98%	0.140%
57	14.751	2014	2017	2025	rVB2	36946	58541	1.74%	0.123%
58	14.928	2039	2047	2051	rBV5	31027	67087	2.00%	0.141%
59	15.075	2066	2072	2075	rBV	1980944	2859976	85.18%	6.008%
60	15.104	2075	2077	2086	rVV	1555537	1807184	53.82%	3.796%
61	15.216	2089	2096	2103	rVB	658928	924948	27.55%	1.943%
62	15.357	2109	2120	2125	rBV2	246149	450880	13.43%	0.947%
63	15.616	2155	2164	2168	rVV	106778	165851	4.94%	0.348%
64	15.657	2168	2171	2174	rVV2	23503	32936	0.98%	0.069%
65	15.698	2174	2178	2180	rVV	33698	49837	1.48%	0.105%
66	15.757	2184	2188	2195	rVB7	31130	70382	2.10%	0.148%
67	15.851	2202	2204	2212	rVB5	33617	61882	1.84%	0.130%
68	15.939	2215	2219	2225	rBV	53001	89798	2.67%	0.189%
69	15.998	2225	2229	2236	rBV3	68969	159435	4.75%	0.335%
70	16.063	2236	2240	2244	rVV2	45755	66487	1.98%	0.140%
71	16.145	2249	2254	2259	rVV3	45226	94204	2.81%	0.198%

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72	16.204	2260	2264	2269	rVV	145295	191623	5.71%	0.403%
73	16.263	2269	2274	2279	rVV3	34293	85688	2.55%	0.180%
74	16.322	2281	2284	2288	rVB	52493	71417	2.13%	0.150%
75	16.398	2294	2297	2302	rVB2	64527	79572	2.37%	0.167%
76	16.445	2302	2305	2310	rVB	28118	35306	1.05%	0.074%
77	16.651	2336	2340	2346	rVV	53664	75943	2.26%	0.160%
78	16.710	2346	2350	2358	rVB2	44285	72979	2.17%	0.153%
79	16.828	2363	2370	2380	rVB	1386458	1902584	56.67%	3.997%
80	16.927	2380	2387	2395	rBV	2153577	2991568	89.10%	6.285%
81	17.028	2398	2404	2413	rBV	2295299	2722709	81.09%	5.720%
82	17.892	2546	2551	2556	rVB3	76988	103808	3.09%	0.218%
83	18.292	2616	2619	2625	rVB	42630	53011	1.58%	0.111%
84	18.574	2663	2667	2677	rBV6	67639	98119	2.92%	0.206%
85	18.745	2692	2696	2701	rVV3	33366	52213	1.56%	0.110%
86	18.963	2730	2733	2737	rVB	37746	44138	1.31%	0.093%
87	19.010	2738	2741	2746	rVB	37745	43749	1.30%	0.092%
88	19.116	2755	2759	2763	rVB2	21297	32552	0.97%	0.068%
89	19.233	2773	2779	2786	rBV	2378159	3357530	100.00%	7.053%
90	19.292	2787	2789	2794	rVB2	29109	36117	1.08%	0.076%
91	20.257	2949	2953	2957	rBV	313161	372511	11.09%	0.783%
92	20.292	2957	2959	2970	rVB4	51882	81429	2.43%	0.171%
93	20.786	3039	3043	3051	rBV	473774	597185	17.79%	1.255%
94	21.033	3080	3085	3095	rBV	1661430	2108654	62.80%	4.430%
95	21.168	3104	3108	3113	rVB	292573	342494	10.20%	0.719%
96	22.080	3261	3263	3267	rVB	130532	125379	3.73%	0.263%
97	22.551	3340	3343	3348	rVB	152269	187379	5.58%	0.394%
98	23.004	3414	3420	3427	rBV	1820048	3035868	90.42%	6.378%
99	23.068	3428	3431	3436	rVV	202245	306989	9.14%	0.645%
100	23.139	3437	3443	3453	rVB	1216898	2096860	62.45%	4.405%

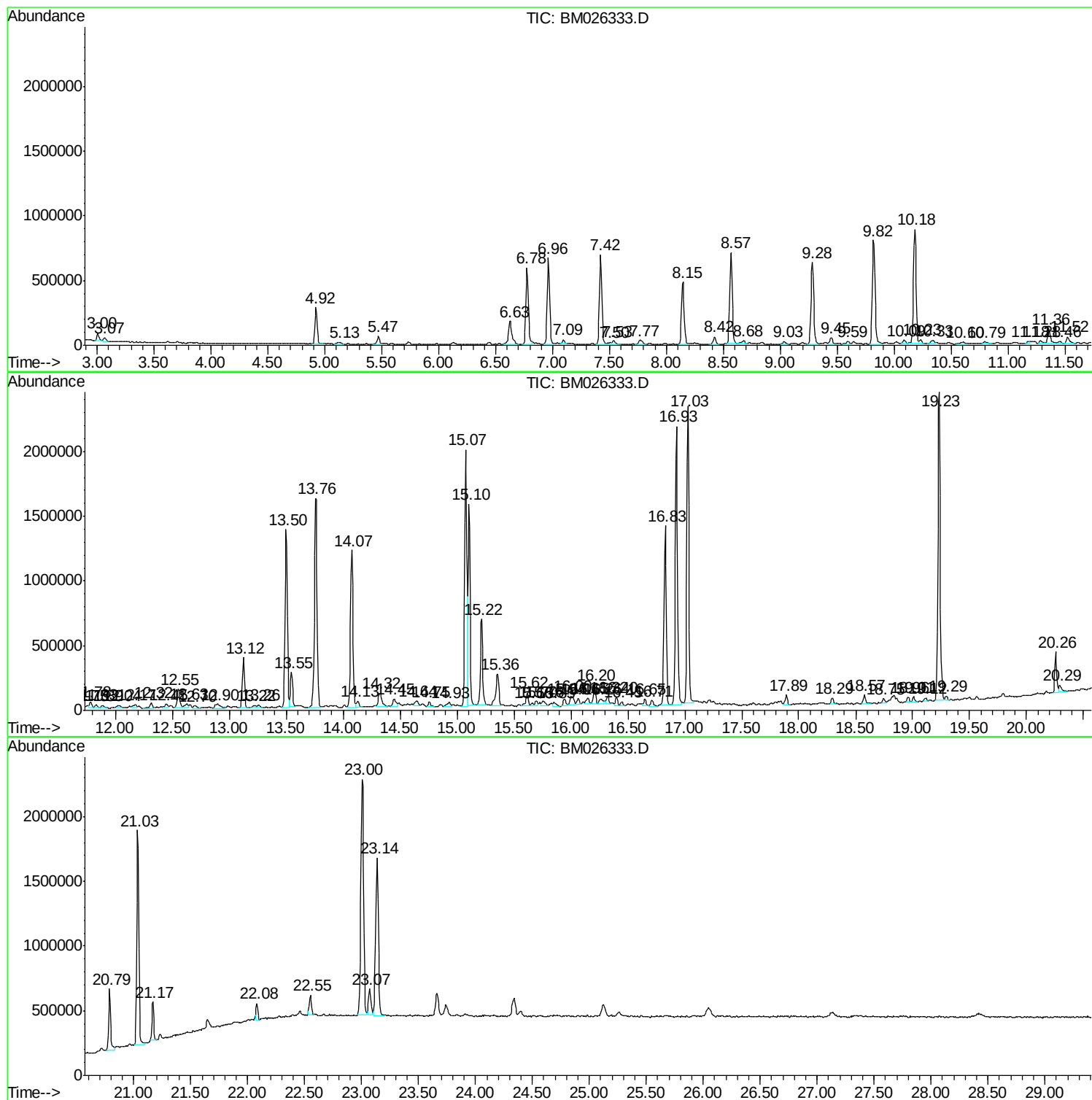
Sum of corrected areas: 47601892

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



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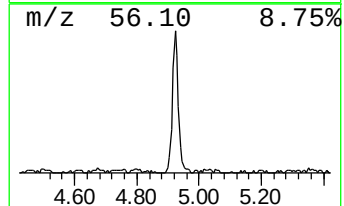
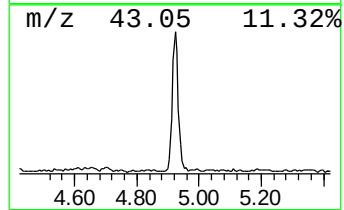
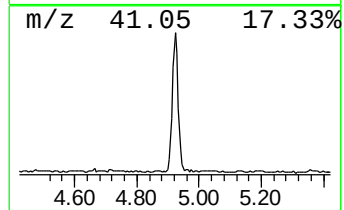
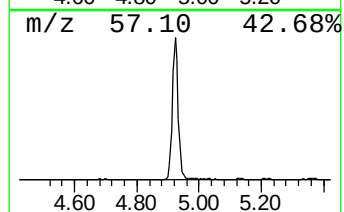
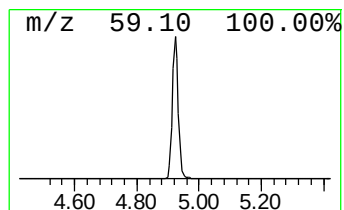
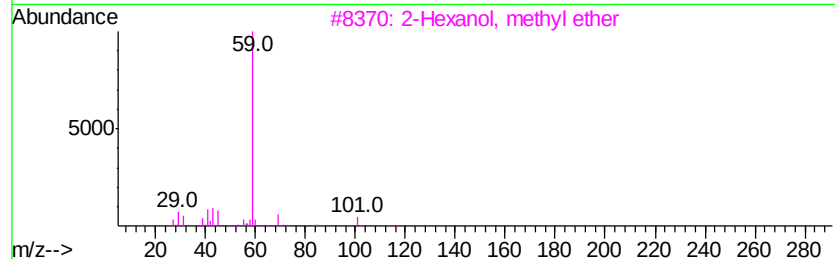
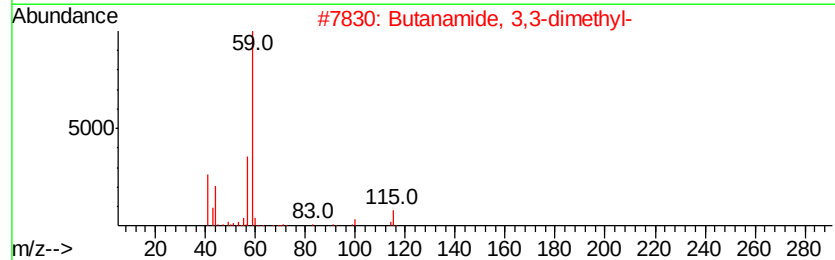
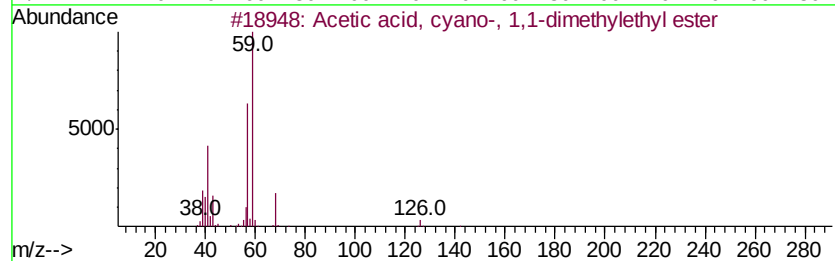
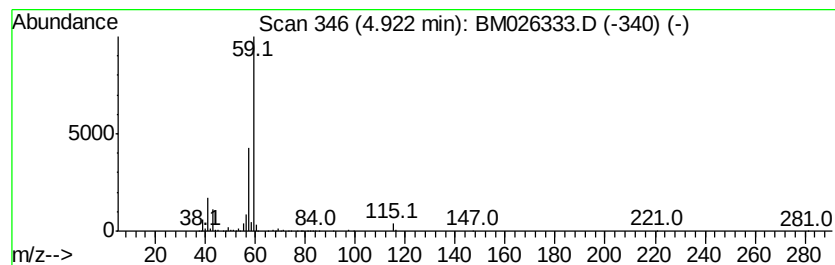
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Acetic acid, cyano-, 1,1-di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	7.33 ng/ul	387049	1,4-Dichlorobenzene-d4	7.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	64
2		Butanamide, 3,3-dimethyl-	115	C6H13NO	000926-04-5	64
3		2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	50
4		Propanoic acid, 2-methoxy-, meth...	118	C5H10O3	017639-76-8	9
5		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9



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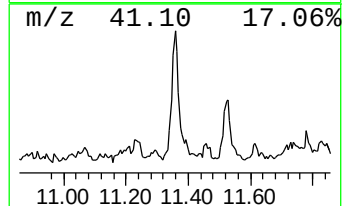
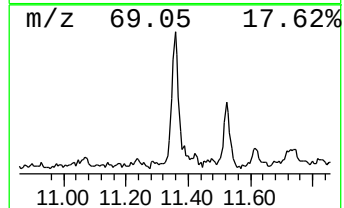
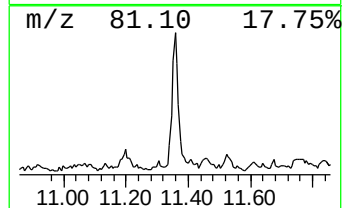
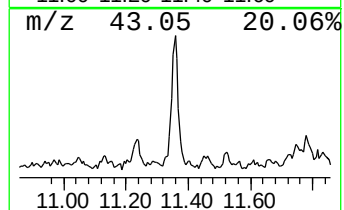
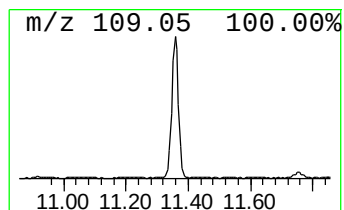
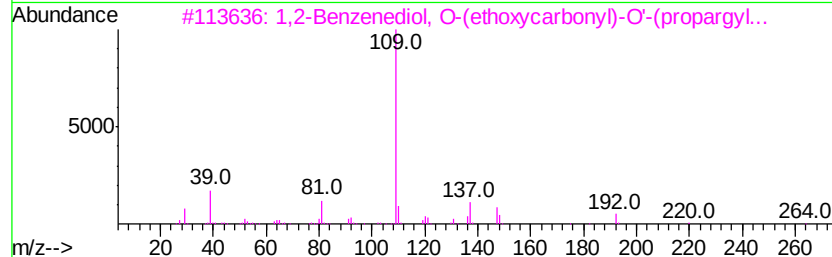
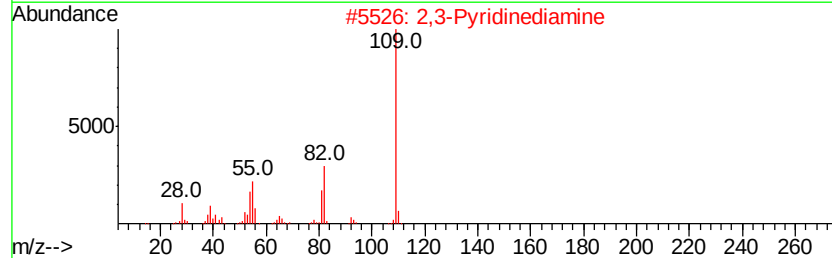
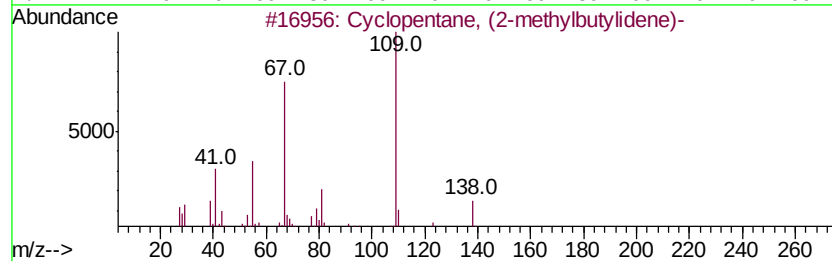
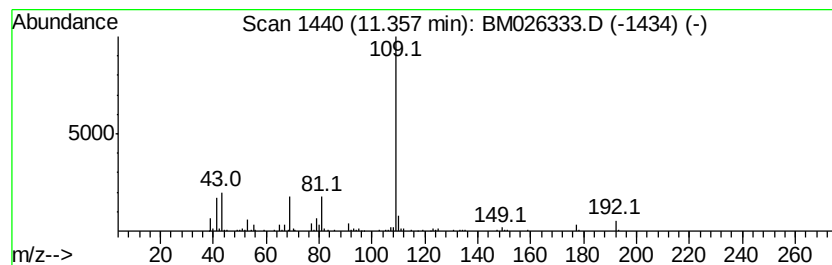
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, (2-methylbuty... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	2.58 ng/ul	184172	Naphthalene-d8	10.18
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentane, (2-methylbutylidene)-	138 C10H18	053366-54-4 59
2		2,3-Pyridinediamine	109 C5H7N3	000452-58-4 59
3		1,2-Benzenediol, O-(ethoxycarbon...	264 C13H12O6	1000329-71-6 50
4		7-(2,6-Dimethyl-hepta-1,5-dienyl...	272 C20H32	1000190-62-8 45
5		(2-Fluorophenyl) methanol, 1-met...	182 C11H15FO	1000374-56-6 42



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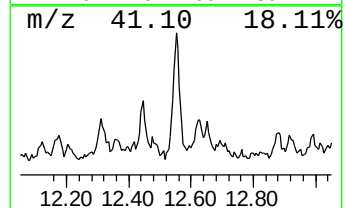
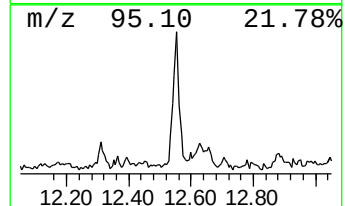
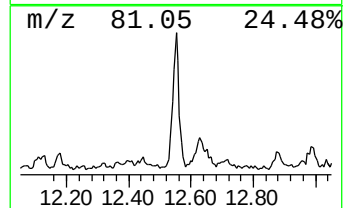
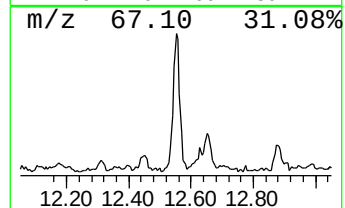
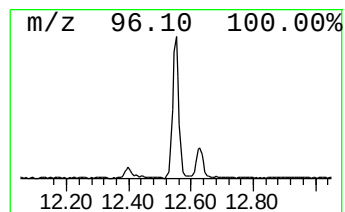
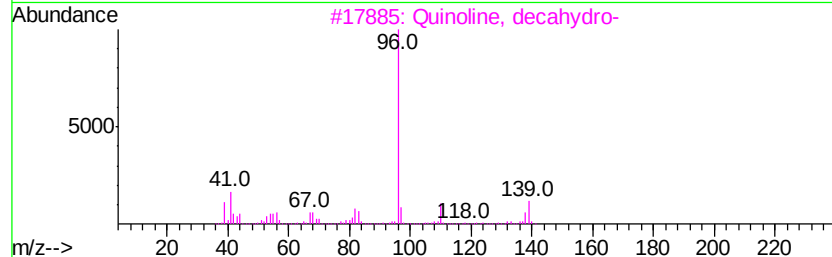
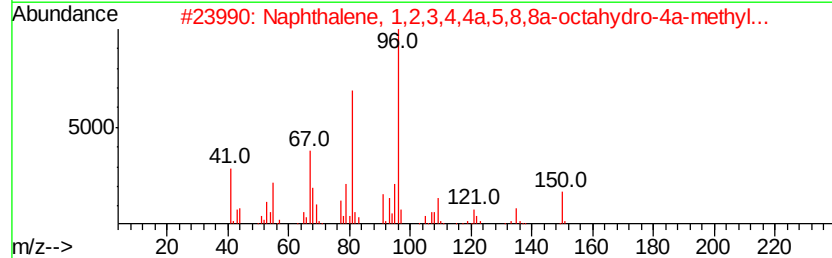
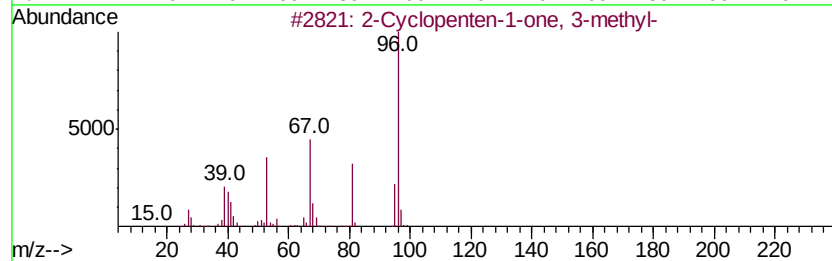
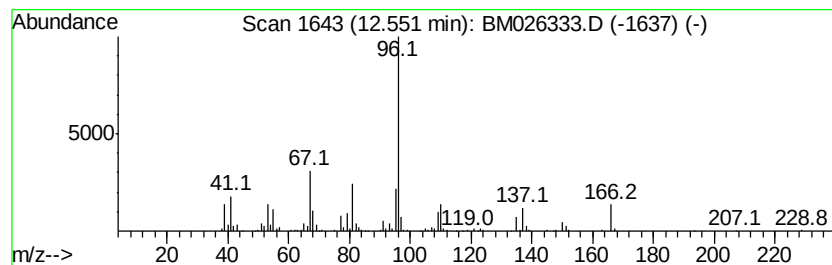
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Peak Number 3 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	2.64 ng/ul	230727	Acenaphthene-d10	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Cyclopenten-1-one, 3-methyl-	96	C6H8O	002758-18-1	58
2		Naphthalene, 1,2,3,4,4a,5,8,8a-o...	150	C11H18	021789-56-0	50
3		Quinoline, decahydro-	139	C9H17N	002051-28-7	46
4		Cyclohexene, 1-decyl-	222	C16H30	062338-41-4	45
5		3,5-Dimethylpyrazole	96	C5H8N2	000067-51-6	43



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Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
Misc :
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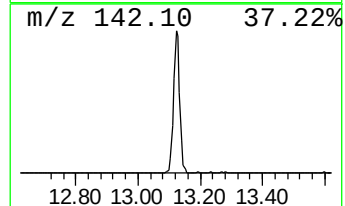
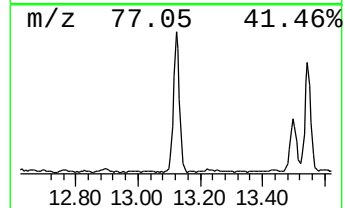
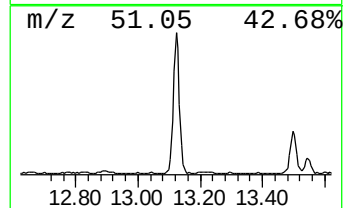
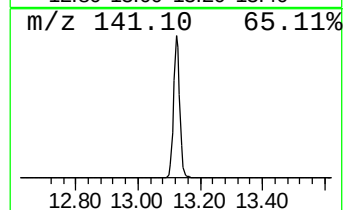
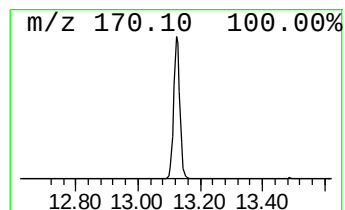
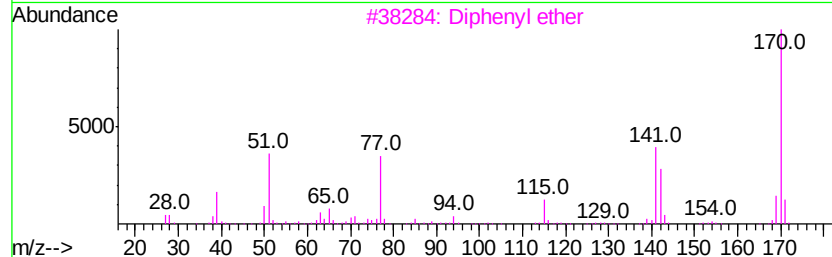
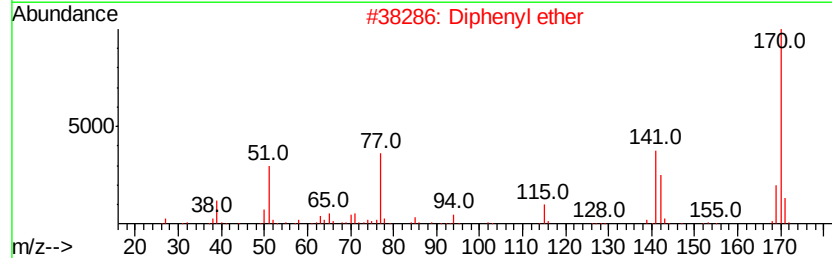
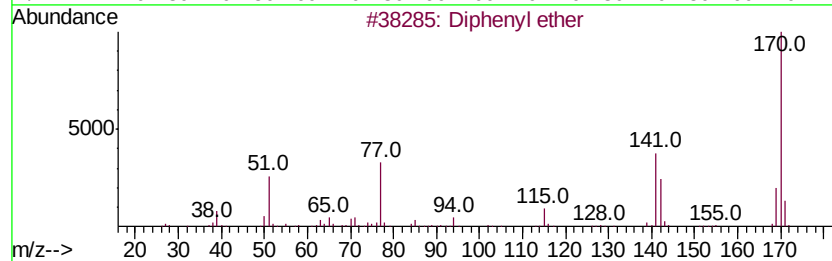
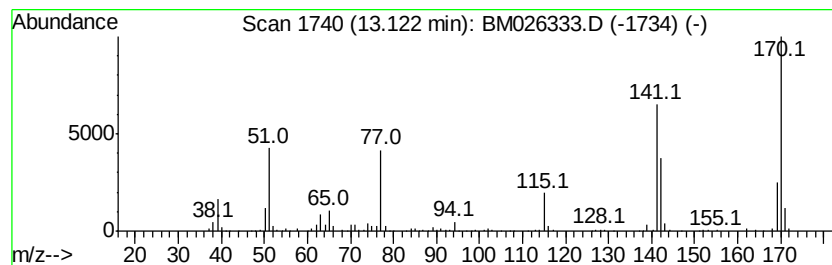
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Diphenyl ether Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	6.19 ng/ul	540363	Acenaphthene-d10	14.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Diphenyl ether	170 C12H10O	000101-84-8	81
2	Diphenyl ether	170 C12H10O	000101-84-8	81
3	Diphenyl ether	170 C12H10O	000101-84-8	74
4	Diphenyl ether	170 C12H10O	000101-84-8	72
5	Spiro[cyclopropane-1,1'(4'H)-nap...	170 C12H10O	033498-24-7	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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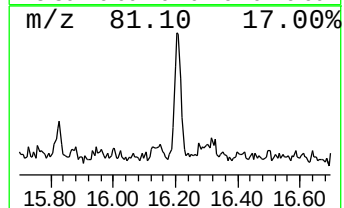
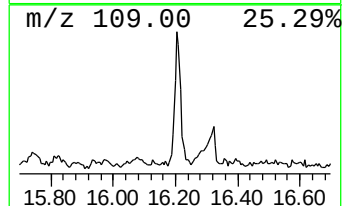
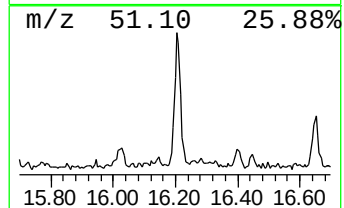
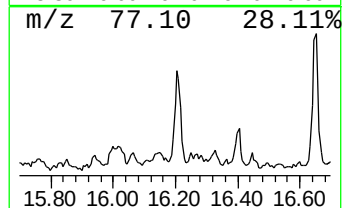
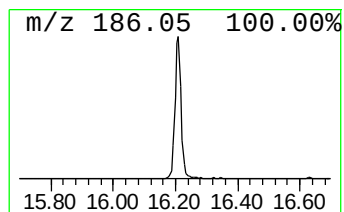
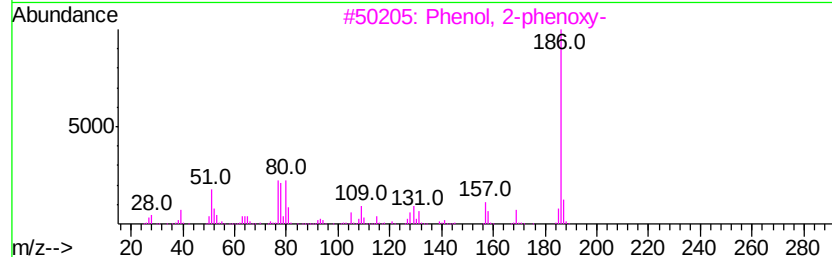
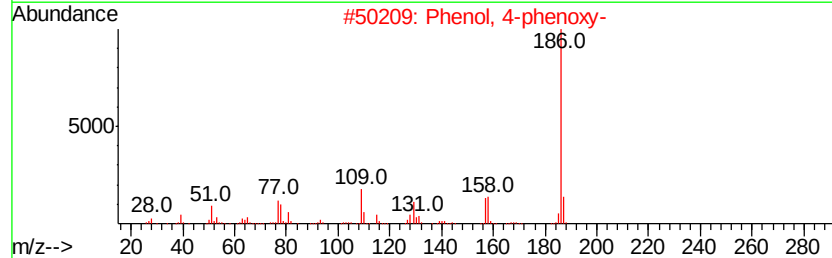
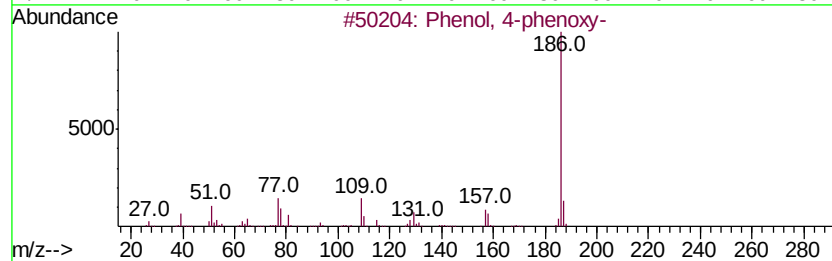
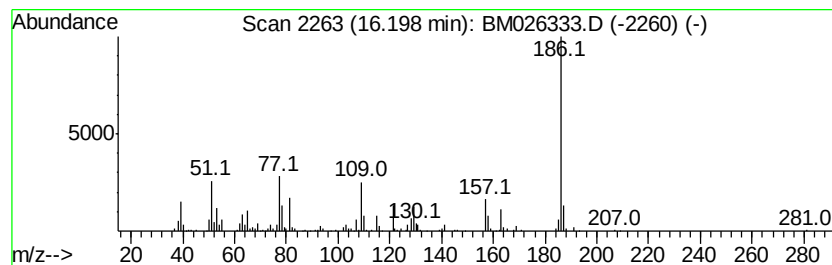
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenol, 4-phenoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	2.01 ng/ul	191623	Phenanthrene-d10	16.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-phenoxy-	186	C12H10O2	000831-82-3	94
2		Phenol, 4-phenoxy-	186	C12H10O2	000831-82-3	89
3		Phenol, 2-phenoxy-	186	C12H10O2	002417-10-9	83
4		Diphenyl sulfide	186	C12H10S	000139-66-2	50
5		[1,1'-Biphenyl]-4,4'-diol	186	C12H10O2	000092-88-6	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

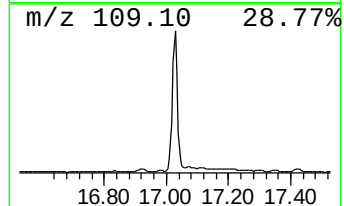
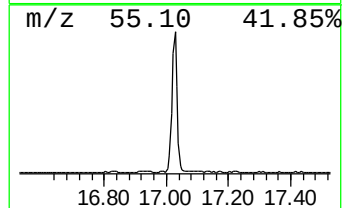
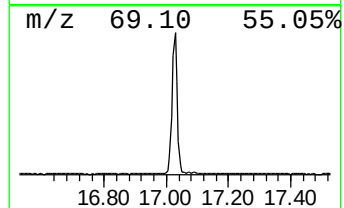
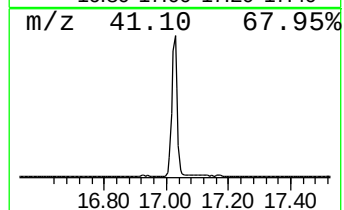
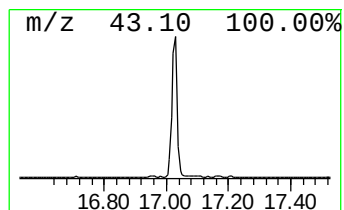
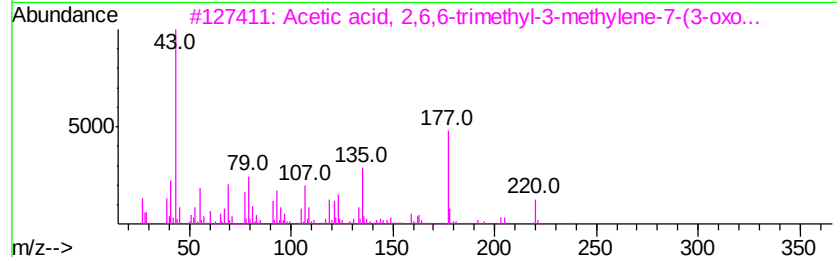
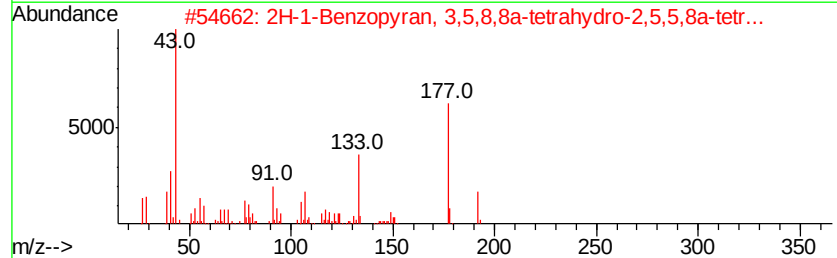
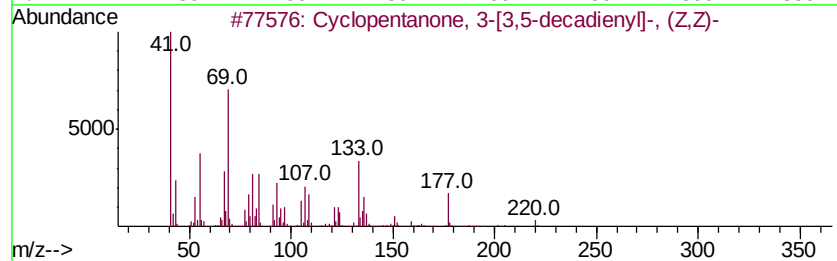
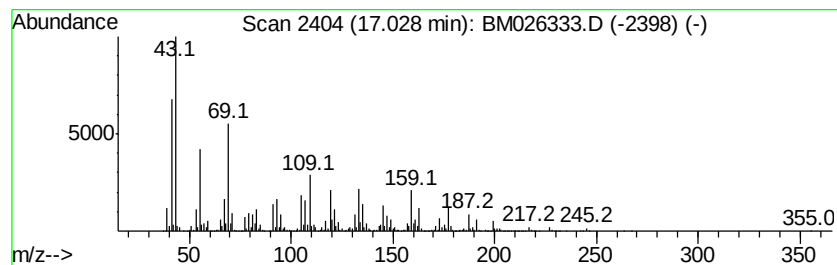
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.03	28.62 ng/ul	2722710	Phenanthrene-d10	16.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopentanone, 3-[3,5-decadieny...	220 C15H24O	1000131-99-8 25
2		2H-1-Benzopyran, 3,5,8,8a-tetra...	192 C13H20O	072468-40-7 22
3		Acetic acid, 2,6,6-trimethyl-3-m...	280 C16H24O4	1000185-41-4 16
4		7-(1,3-Dimethylbuta-1,3-dienyl)-...	234 C15H22O2	1000190-22-7 16
5		Bicyclo[4.3.0]nonan-1-ol, 7,9-bi...	206 C14H22O	125257-63-8 14



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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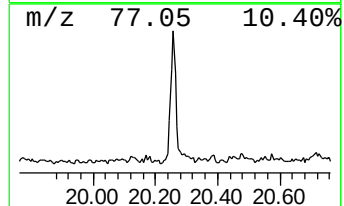
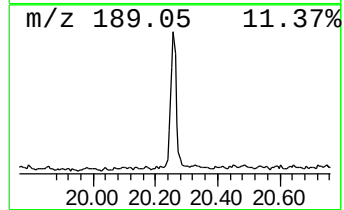
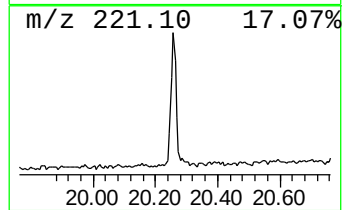
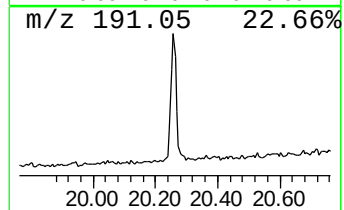
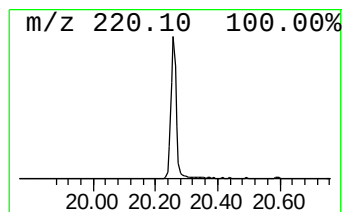
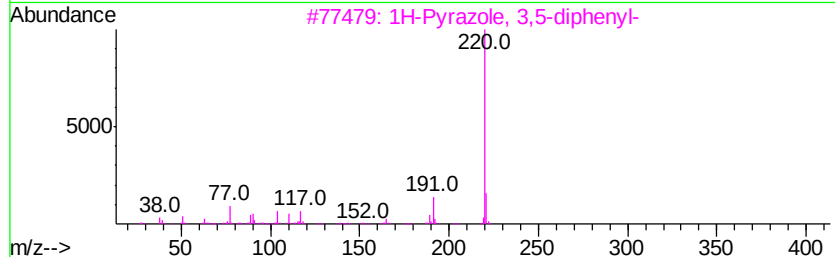
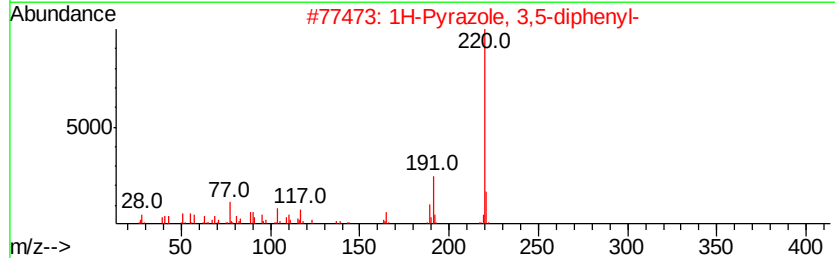
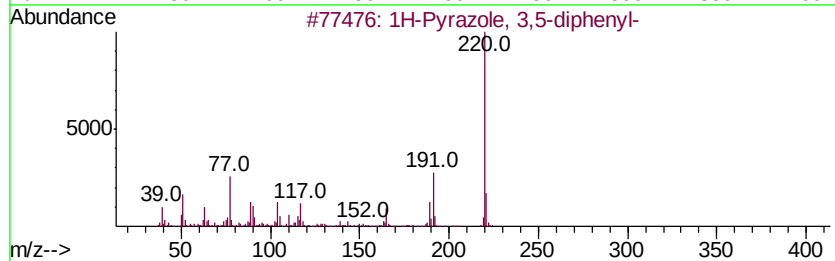
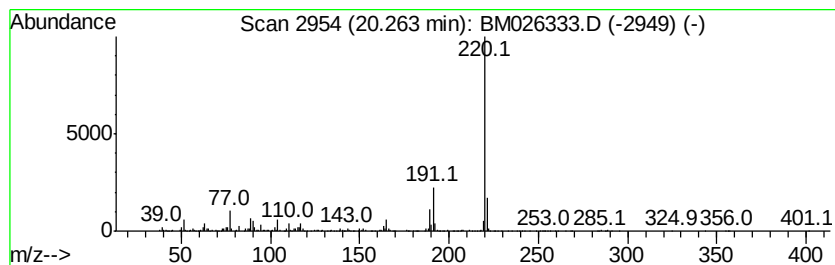
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1H-Pyrazole, 3,5-diphenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.26	3.53 ng/ul	372511	Chrysene-d12	21.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	97
2		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	95
3		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	91
4		1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	90
5		Naphthalene, 2-phenoxy-	220	C16H12O	019420-29-2	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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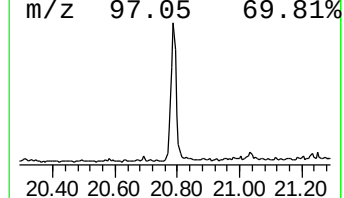
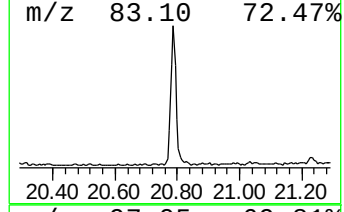
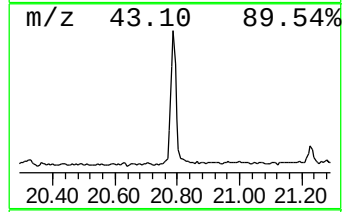
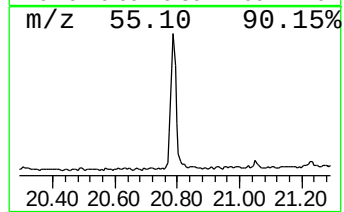
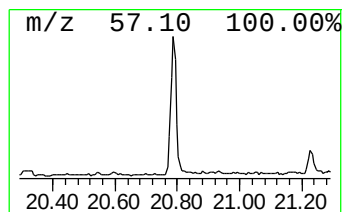
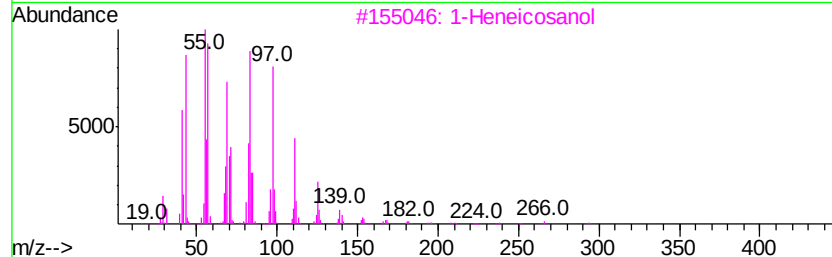
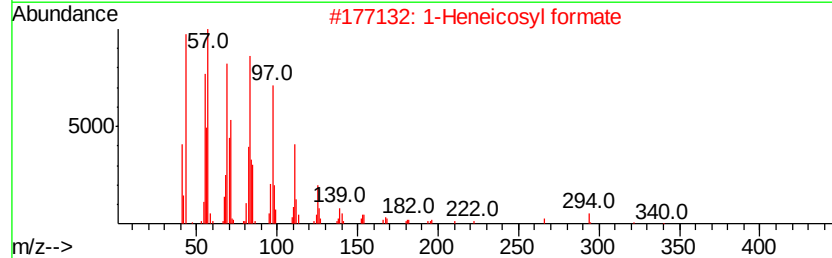
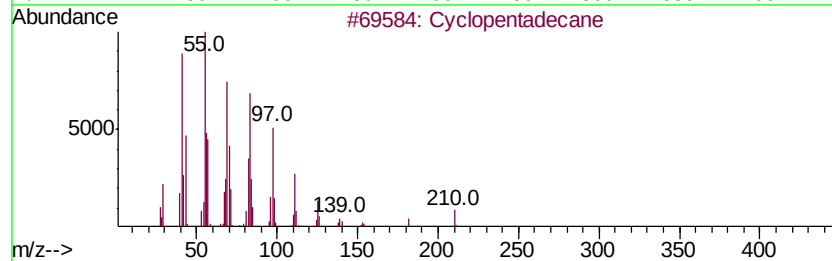
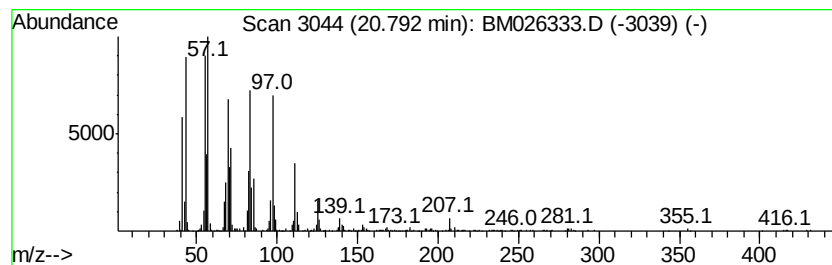
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic20.79 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	5.66 ng/ul	597185	Chrysene-d12	21.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	97
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		Cyclotetracosane	336	C24H48	000297-03-0	93
5		1-Heptacosanol	396	C27H56O	002004-39-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
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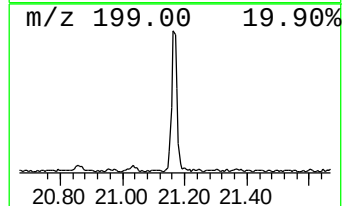
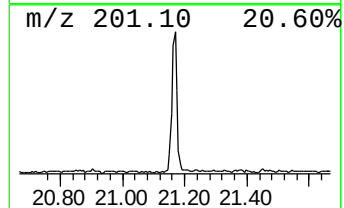
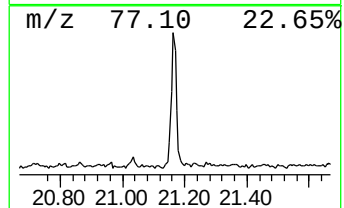
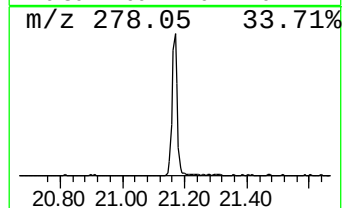
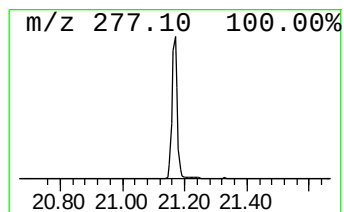
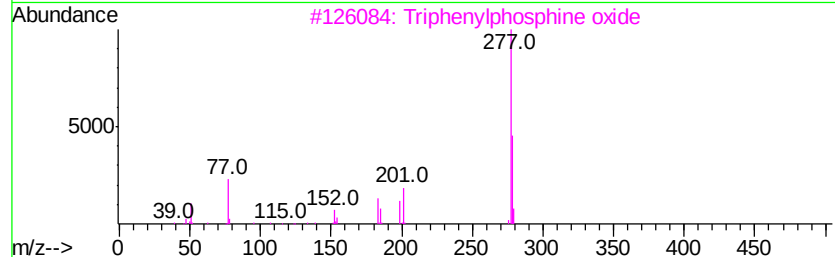
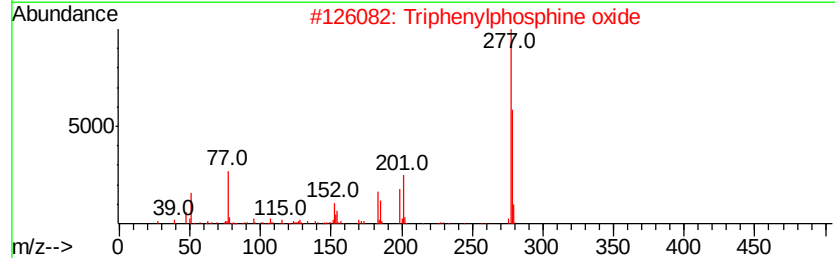
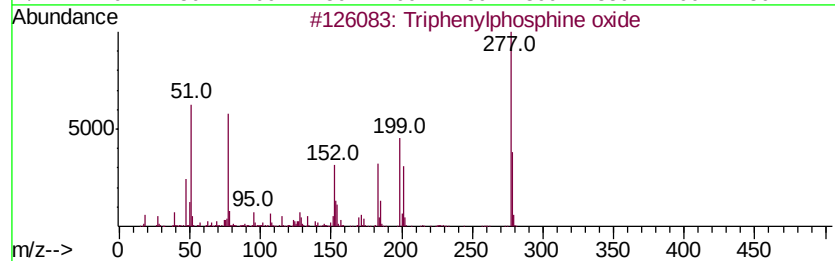
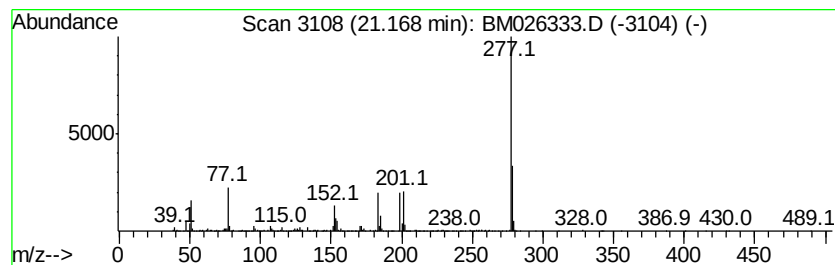
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Triphenylphosphine oxide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.17	3.25 ng/ul	342494	Chrysene-d12		21.03		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	64
3			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	50
4			(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	47
5			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
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Sample : L2884-01
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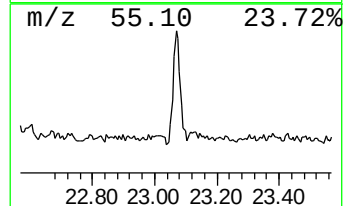
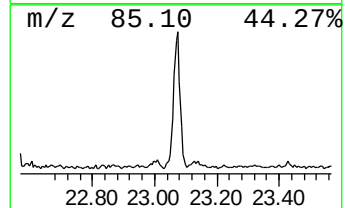
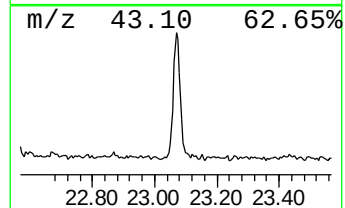
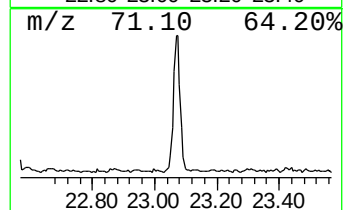
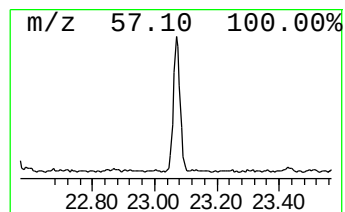
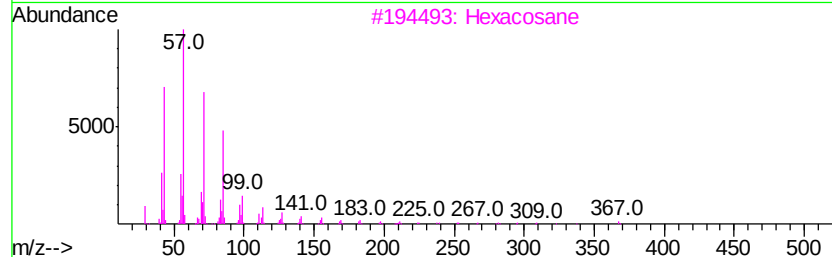
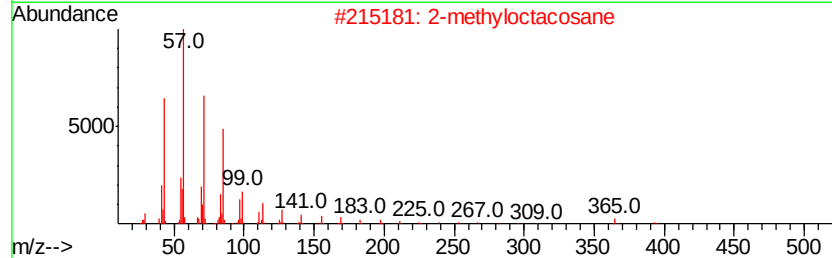
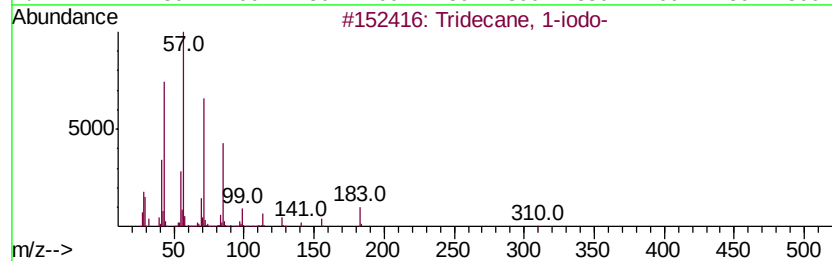
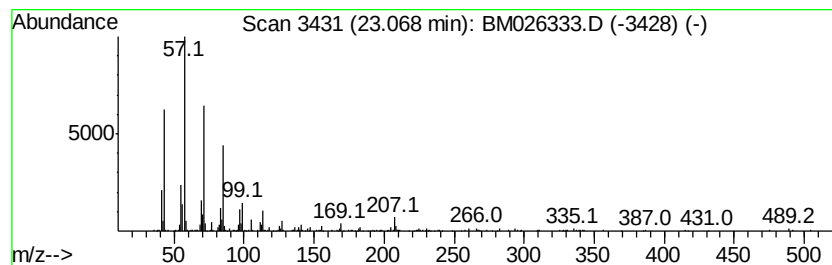
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Tridecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.07	2.93 ng/ul	306989	Perylene-d12	23.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Hexadecane	226	C16H34	000544-76-3	89
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM061020\
Data File : BM026333.D
Acq On : 10 Jun 2020 17:01
Operator : JU/CG
Sample : L2884-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFXD7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Acetic acid, cyan...	4.92	7.3	ng/ul	387049	1	7.42	1055450	20.0
Cyclopentane, (2-...	11.36	2.6	ng/ul	184172	2	10.18	1426710	20.0
2-Cyclopenten-1-o...	12.55	2.6	ng/ul	230727	3	14.07	1746840	20.0
Diphenyl ether	13.12	6.2	ng/ul	540363	3	14.07	1746840	20.0
Phenol, 4-phenoxy-	16.20	2.0	ng/ul	191623	4	16.83	1902580	20.0
unknown-01	17.03	28.6	ng/ul	2722710	4	16.83	1902580	20.0
1H-Pyrazole, 3,5-...	20.26	3.5	ng/ul	372511	5	21.03	2108650	20.0
(DEL) Alkane: Cyc...	20.79	5.7	ng/ul	597185	5	21.03	2108650	20.0
Triphenylphosphin...	21.17	3.3	ng/ul	342494	5	21.03	2108650	20.0
Tridecane, 1-iodo-	23.07	2.9	ng/ul	306989	6	23.14	2096860	20.0