

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
 Data File : BM026977.D
 Acq On : 03 Aug 2020 12:44
 Operator : JU/CG
 Sample : L3424-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0S60

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-BM072720MA.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	2.766	3	5	11	rVB3	10146	12042	0.39%	0.025%
2	2.902	23	28	36	rBV	136238	191429	6.19%	0.393%
3	2.972	36	40	50	rVB	294828	331700	10.72%	0.681%
4	3.219	79	82	84	rBV2	8943	10160	0.33%	0.021%
5	3.249	84	87	94	rVB	27616	34720	1.12%	0.071%
6	4.825	350	355	365	rBV	157326	210616	6.81%	0.433%
7	6.837	690	697	706	rVB	261444	398527	12.88%	0.818%
8	7.001	717	725	732	rBV	180800	284565	9.20%	0.584%
9	7.195	751	758	765	rBV	253783	395167	12.77%	0.811%
10	7.301	772	776	784	rVB4	2974	6676	0.22%	0.014%
11	7.660	830	837	843	rVB	266420	407223	13.16%	0.836%
12	8.195	919	928	934	rBV2	97402	227020	7.34%	0.466%
13	8.248	934	937	940	rVB3	9457	12417	0.40%	0.025%
14	8.360	949	956	965	rBV	295108	456049	14.74%	0.937%
15	8.801	1024	1031	1035	rBV	243092	411303	13.29%	0.845%
16	8.842	1035	1038	1047	rVB	187598	278057	8.99%	0.571%
17	9.519	1146	1153	1156	rBV	185305	331013	10.70%	0.680%
18	9.548	1156	1158	1166	rVB	200017	280352	9.06%	0.576%
19	9.854	1204	1210	1216	rBV	153714	227556	7.36%	0.467%
20	10.054	1237	1244	1248	rBV	322145	642527	20.77%	1.319%
21	10.436	1301	1309	1313	rBV	334519	572896	18.52%	1.176%
22	10.484	1313	1317	1324	rVB	181039	287389	9.29%	0.590%
23	10.566	1324	1331	1344	rVB	270901	472286	15.27%	0.970%
24	11.719	1520	1527	1537	rBV	833274	1262844	40.82%	2.593%
25	11.972	1566	1570	1575	rBV4	6353	10897	0.35%	0.022%
26	12.025	1575	1579	1585	rVV2	4254	6729	0.22%	0.014%
27	12.101	1585	1592	1601	rVB	561115	848465	27.43%	1.742%
28	12.236	1608	1615	1621	rBV5	12649	17672	0.57%	0.036%
29	12.325	1621	1630	1638	rVB	455369	721142	23.31%	1.481%
30	12.466	1648	1654	1660	rBV	337102	499430	16.14%	1.026%
31	12.719	1690	1697	1705	rVB	255895	392751	12.70%	0.807%
32	12.913	1726	1730	1735	rVB2	5298	7459	0.24%	0.015%
33	13.160	1764	1772	1778	rBV	437884	655966	21.20%	1.347%
34	13.213	1778	1781	1785	rVB2	9644	12129	0.39%	0.025%

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35	13.707	1858	1865	1869	rBV	679346	889148	28.74%	1.826%
36	13.754	1869	1873	1882	rVB	153623	204912	6.62%	0.421%
37	13.860	1885	1891	1897	rVB	195291	260304	8.41%	0.535%
38	13.983	1901	1912	1915	rBV	653451	1080354	34.92%	2.219%
39	14.295	1958	1965	1970	rBV	504702	746562	24.13%	1.533%
40	14.354	1970	1975	1982	rVB	353809	513024	16.58%	1.054%
41	14.483	1991	1997	2007	rBV	360107	501975	16.23%	1.031%
42	14.654	2020	2026	2035	rVV	509404	693648	22.42%	1.424%
43	14.919	2065	2071	2081	rBV	677685	943936	30.51%	1.938%
44	15.124	2096	2106	2110	rBV2	16633	24508	0.79%	0.050%
45	15.177	2112	2115	2120	rVB3	8602	11519	0.37%	0.024%
46	15.289	2127	2134	2138	rBV	876589	1232576	39.84%	2.531%
47	15.342	2138	2143	2148	rVV	981664	1302881	42.11%	2.676%
48	15.418	2148	2156	2167	rVV3	1210177	1840945	59.51%	3.780%
49	15.524	2169	2174	2180	rVB5	11365	20896	0.68%	0.043%
50	15.660	2191	2197	2202	rBV8	5148	9103	0.29%	0.019%
51	15.936	2240	2244	2249	rBV4	10879	15074	0.49%	0.031%
52	16.236	2285	2295	2302	rVB	673602	906302	29.30%	1.861%
53	16.307	2302	2307	2309	rBV3	14248	19285	0.62%	0.040%
54	16.348	2309	2314	2320	rVV	710977	965783	31.22%	1.983%
55	16.460	2329	2333	2337	rVB5	5534	9946	0.32%	0.020%
56	16.689	2367	2372	2375	rBV6	3894	6570	0.21%	0.013%
57	16.830	2392	2396	2400	rVV4	6028	10769	0.35%	0.022%
58	16.989	2416	2423	2425	rBV5	6026	8746	0.28%	0.018%
59	17.036	2425	2431	2434	rVV	638487	901050	29.13%	1.850%
60	17.077	2434	2438	2442	rVV	556019	729324	23.57%	1.498%
61	17.130	2442	2447	2451	rVV2	1054723	1607855	51.97%	3.302%
62	17.165	2451	2453	2459	rVB	634844	730178	23.60%	1.499%
63	17.436	2495	2499	2504	rVB5	5371	7508	0.24%	0.015%
64	17.507	2506	2511	2518	rVV	746028	1022393	33.05%	2.100%
65	17.712	2544	2546	2551	rVB5	6245	8528	0.28%	0.018%
66	17.854	2569	2570	2575	rVB5	5630	6958	0.22%	0.014%
67	17.954	2582	2587	2593	rVV	107846	154189	4.98%	0.317%
68	18.018	2595	2598	2602	rVB2	10675	14093	0.46%	0.029%
69	18.448	2666	2671	2680	rVB3	37098	87177	2.82%	0.179%
70	18.554	2686	2689	2693	rBV4	11683	16727	0.54%	0.034%
71	18.730	2712	2719	2723	rVV	208904	268366	8.67%	0.551%

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72	18.765	2724	2725	2730	rVV3	12613	15729	0.51%	0.032%
73	18.818	2730	2734	2740	rVV	53705	68458	2.21%	0.141%
74	18.889	2742	2746	2754	rVB3	18578	32130	1.04%	0.066%
75	19.095	2773	2781	2786	rBV	1656680	2110461	68.22%	4.334%
76	19.236	2800	2805	2810	rBV3	39447	53616	1.73%	0.110%
77	19.318	2815	2819	2824	rVV3	21201	39992	1.29%	0.082%
78	19.365	2824	2827	2832	rVV	32805	47004	1.52%	0.097%
79	19.430	2832	2838	2841	rVV	1443767	2072699	67.00%	4.256%
80	19.454	2841	2842	2851	rVB	672566	646383	20.89%	1.327%
81	19.901	2916	2918	2924	rVB3	8299	10253	0.33%	0.021%
82	19.971	2927	2930	2933	rVB5	12394	14574	0.47%	0.030%
83	20.112	2952	2954	2958	rBV5	7495	8673	0.28%	0.018%
84	20.383	2997	3000	3003	rVB3	12298	14126	0.46%	0.029%
85	20.595	3032	3036	3040	rBV	47173	63079	2.04%	0.130%
86	20.959	3094	3098	3104	rBV	432118	510048	16.49%	1.047%
87	21.218	3135	3142	3146	rBV2	1087729	2110478	68.22%	4.334%
88	21.259	3146	3149	3154	rVB	768100	892389	28.85%	1.833%
89	21.424	3170	3177	3178	rBV4	124393	249684	8.07%	0.513%
90	21.842	3245	3248	3254	rVB4	53560	75212	2.43%	0.154%
91	21.983	3270	3272	3278	rVB	59492	70117	2.27%	0.144%
92	22.771	3400	3406	3411	rBV	1833837	3093677	100.00%	6.353%
93	22.818	3411	3414	3421	rVB	605919	879763	28.44%	1.807%
94	23.289	3488	3494	3499	rBV	1058764	1923519	62.18%	3.950%
95	23.336	3499	3502	3511	rVV	386289	702084	22.69%	1.442%
96	23.430	3512	3518	3530	rVB2	587547	1098451	35.51%	2.256%
97	24.006	3612	3616	3622	rVB3	77890	137822	4.45%	0.283%
98	24.077	3624	3628	3635	rVB4	75604	147246	4.76%	0.302%
99	25.641	3884	3894	3905	rVB3	642464	1933778	62.51%	3.971%
100	26.300	3998	4006	4023	rVB	325688	984610	31.83%	2.022%

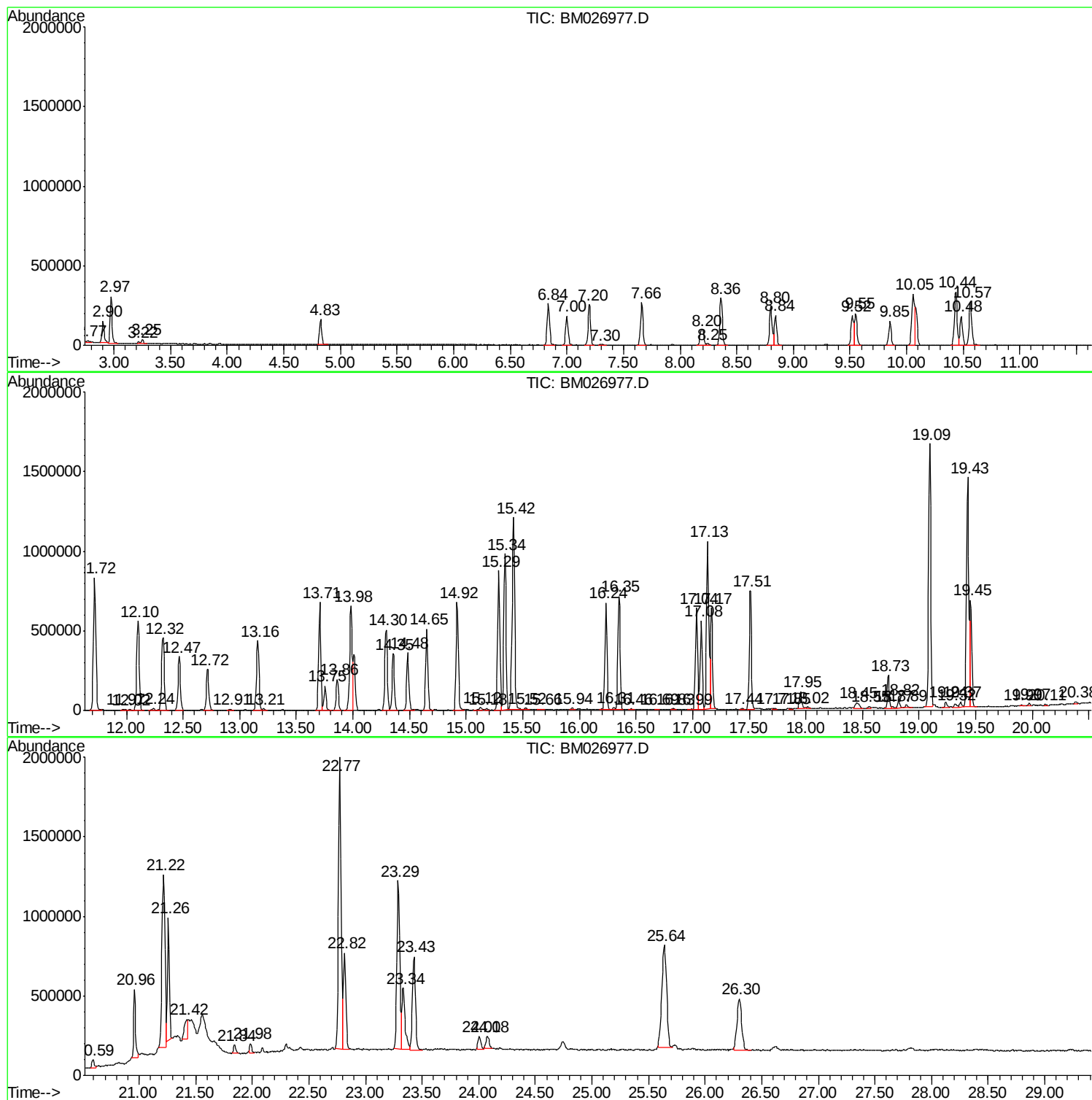
Sum of corrected areas: 48696341

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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



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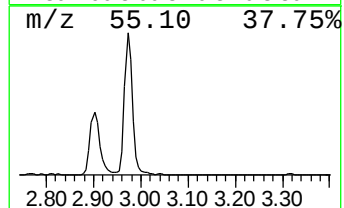
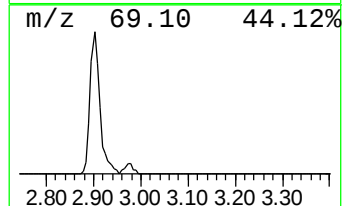
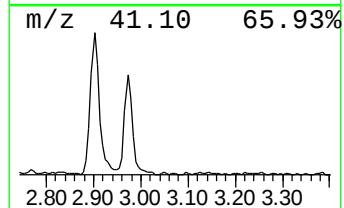
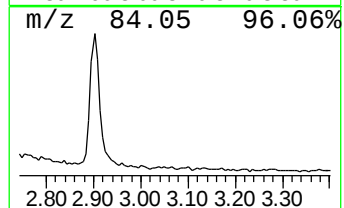
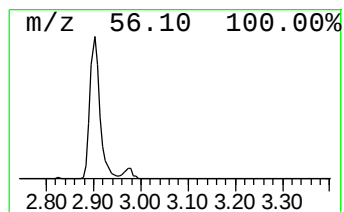
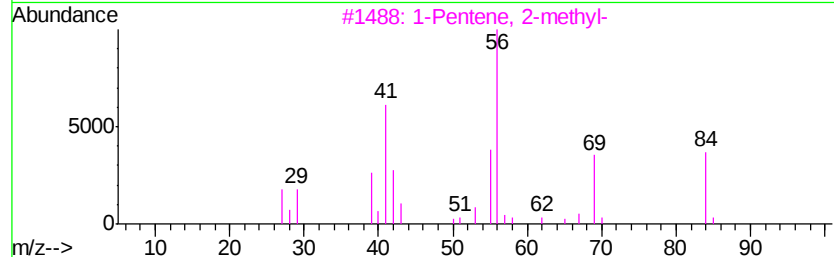
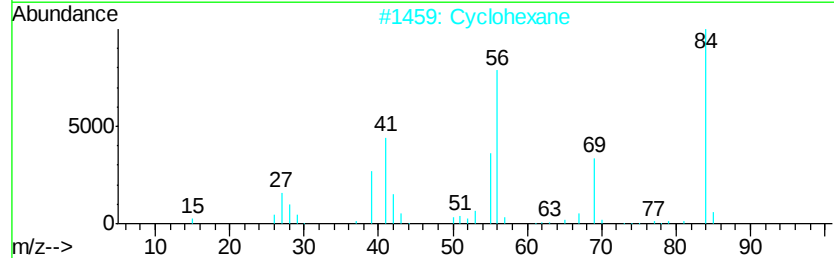
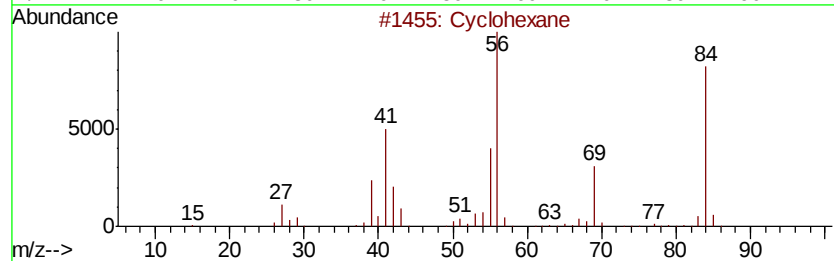
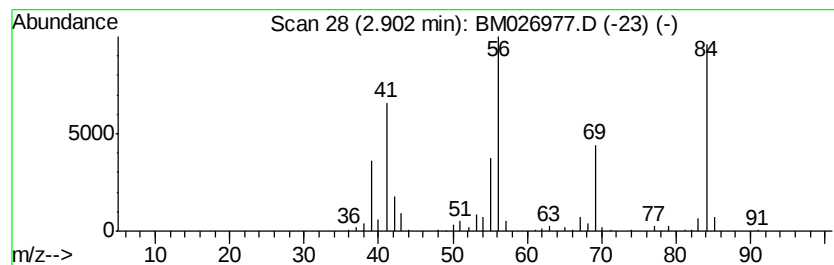
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic2.90 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.90	9.40 ng/ul	191429	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	90
2			Cyclohexane	84	C6H12	000110-82-7	90
3			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	86
4			Cyclohexane	84	C6H12	000110-82-7	83
5			Cyclohexane	84	C6H12	000110-82-7	83



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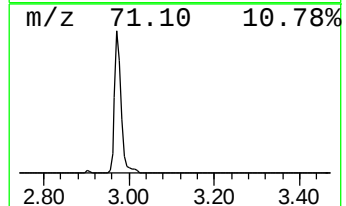
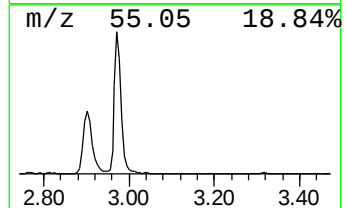
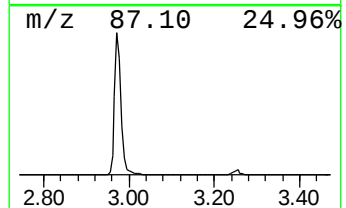
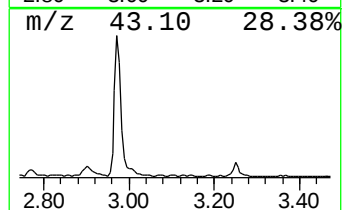
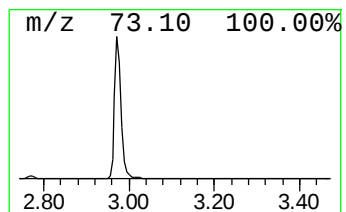
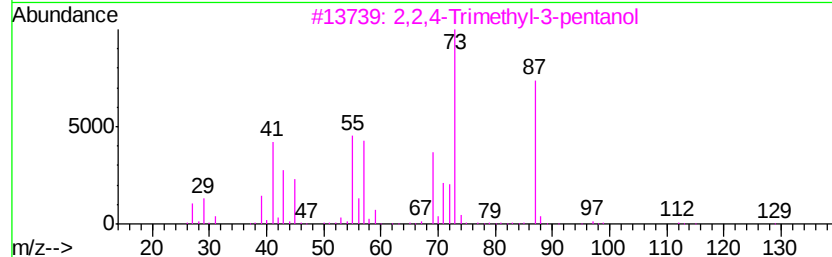
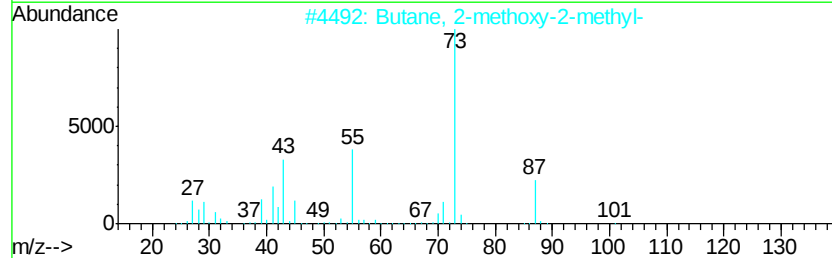
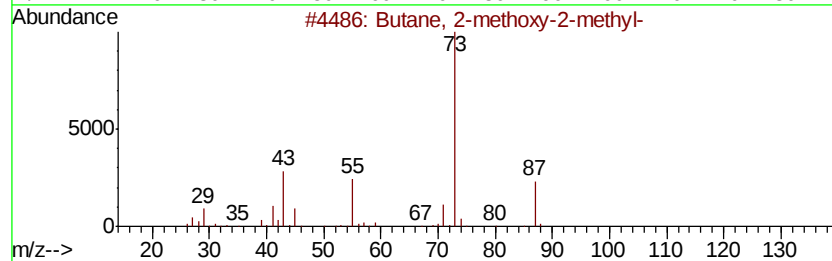
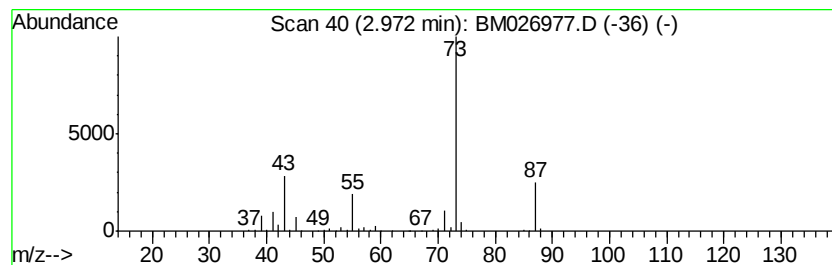
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.97	16.29 ng/ul	331700	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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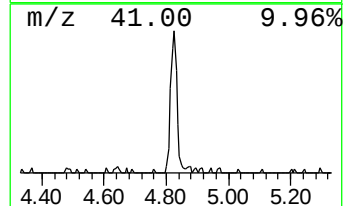
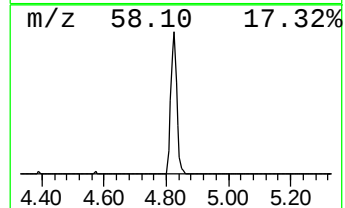
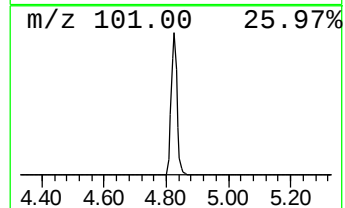
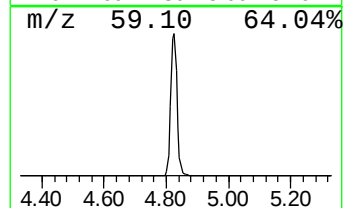
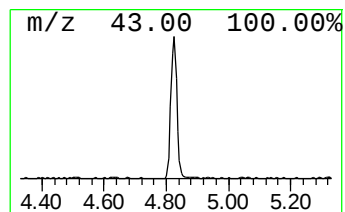
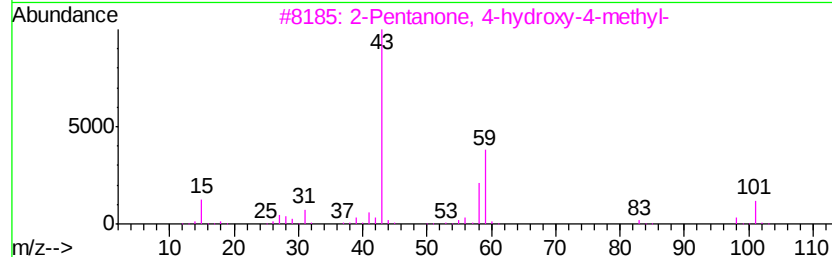
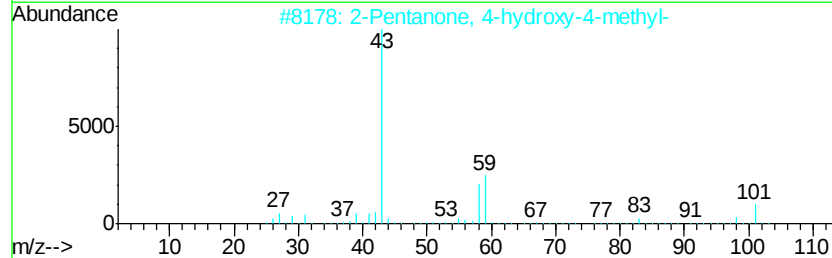
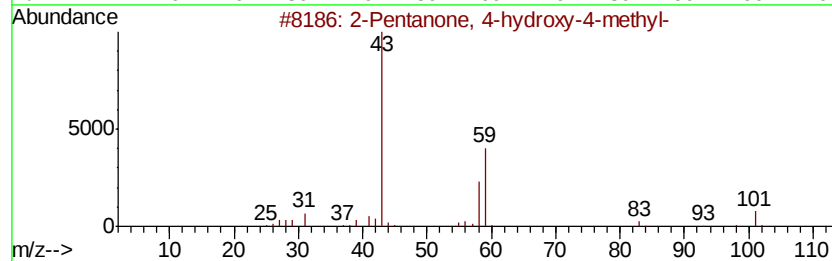
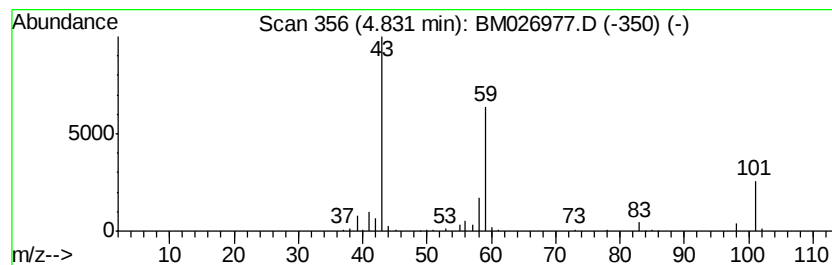
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	10.34 ng/ul	210616	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
4		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
5		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
Operator : JU/CG
Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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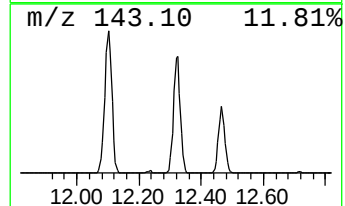
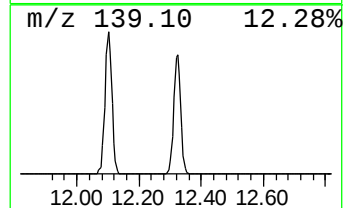
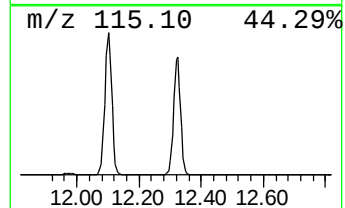
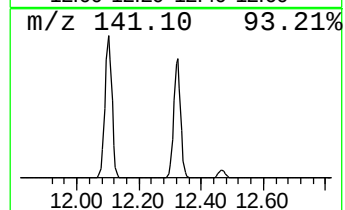
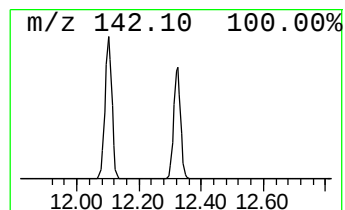
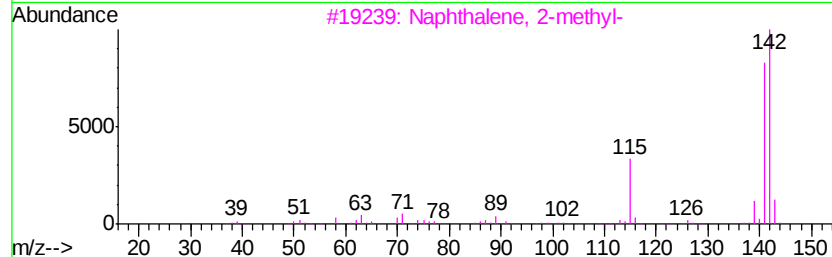
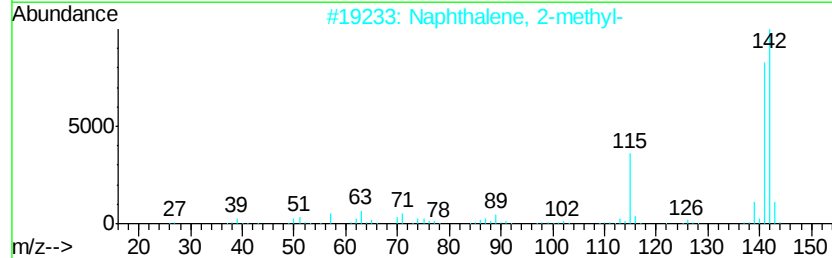
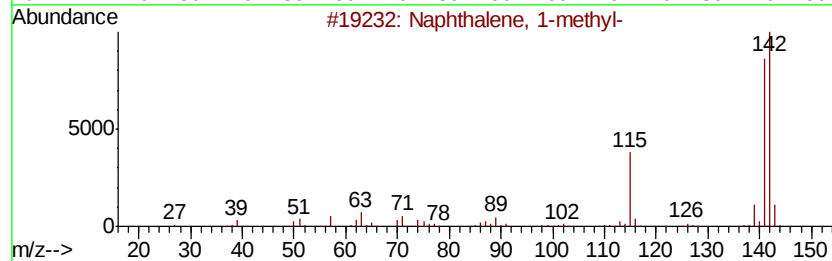
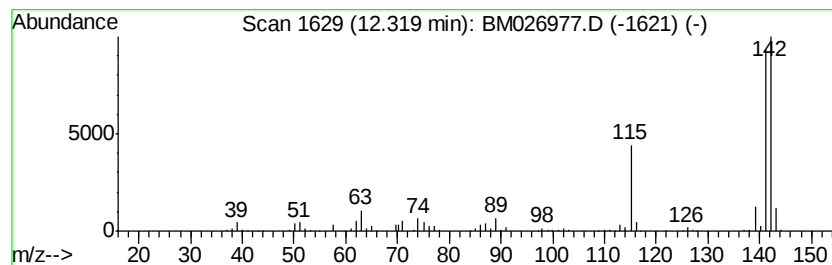
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	25.18 ng/ul	721142	Naphthalene-d8	10.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
Operator : JU/CG
Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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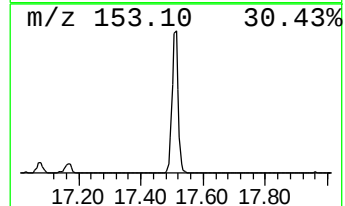
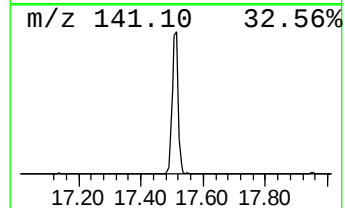
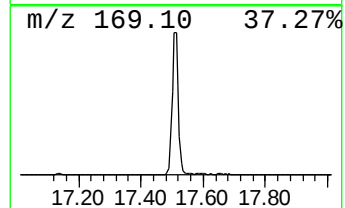
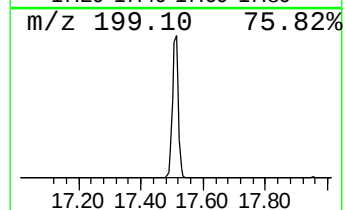
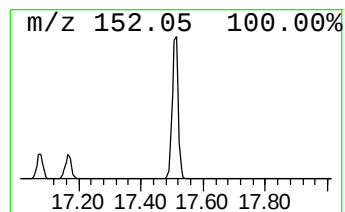
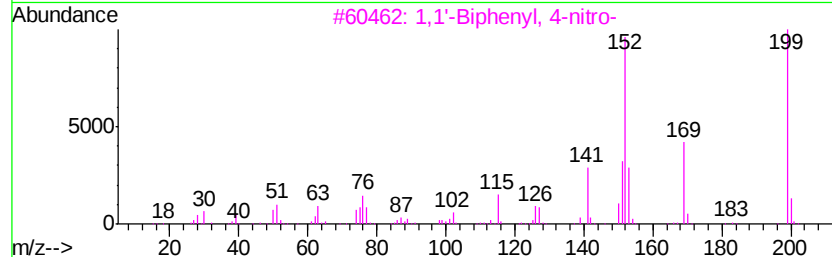
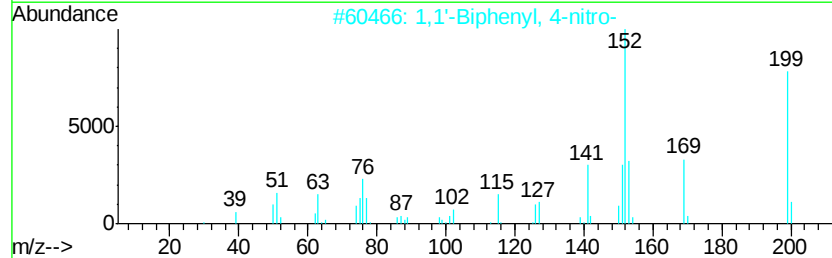
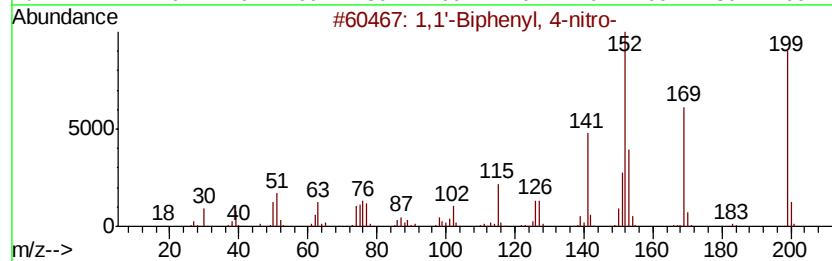
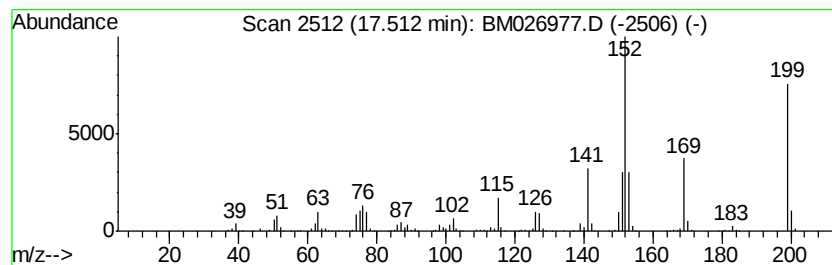
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.51	22.69 ng/ul	1022390	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
2		1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3		1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
4		Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	94
5		1,1'-Biphenyl, 3-nitro-	199	C12H9NO2	002113-58-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
Operator : JU/CG
Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

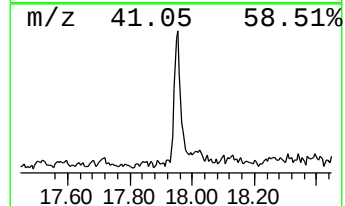
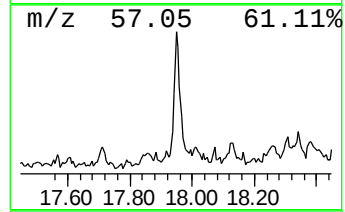
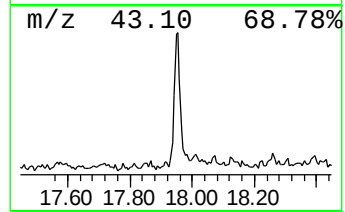
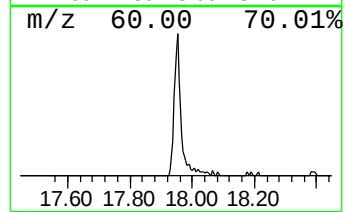
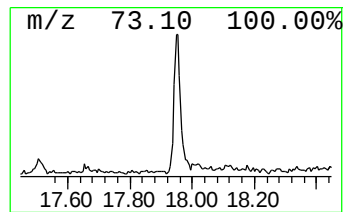
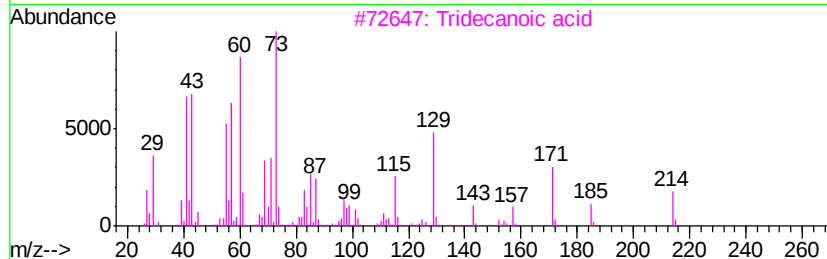
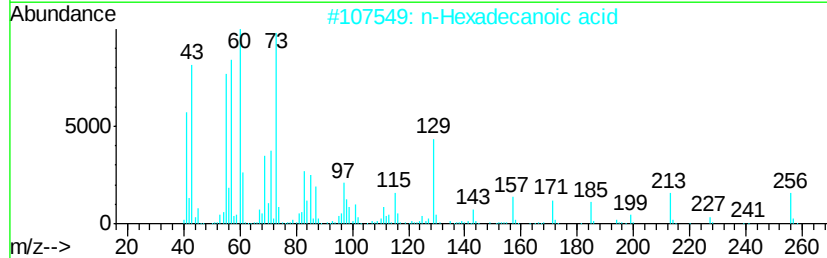
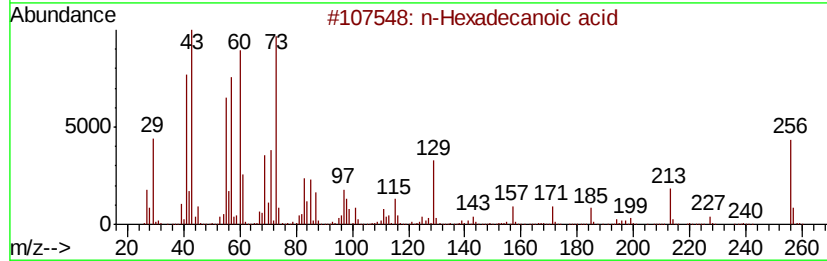
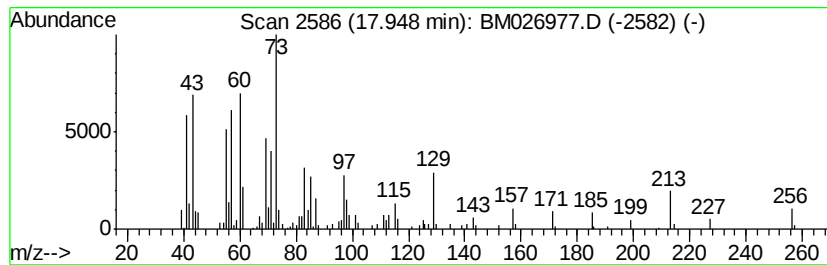
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.95	3.42 ng/ul	154189	Phenanthrene-d10	17.04		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94	
3	Tridecanoic acid	214	C13H26O2	000638-53-9	94	
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87	
5	Tridecanoic acid	214	C13H26O2	000638-53-9	83	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
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Acq On : 03 Aug 2020 12:44
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Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

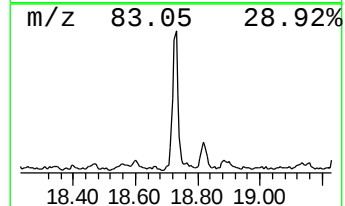
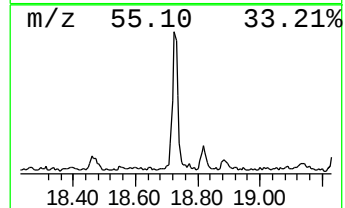
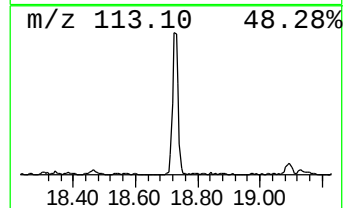
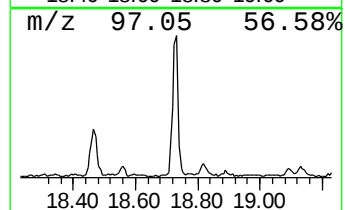
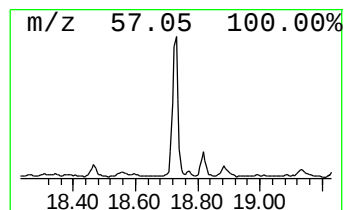
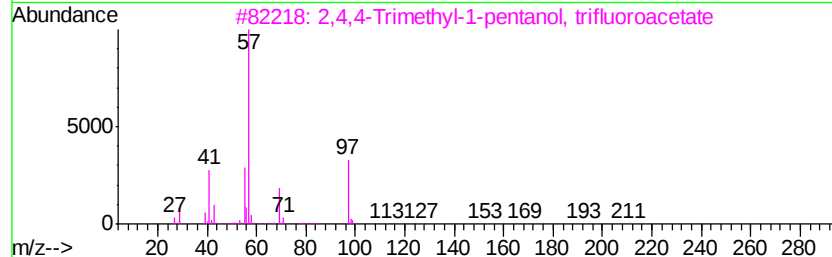
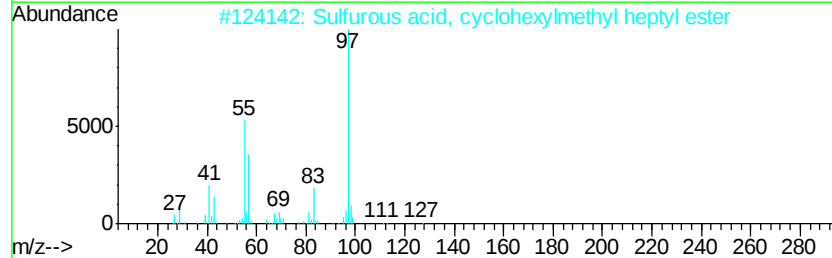
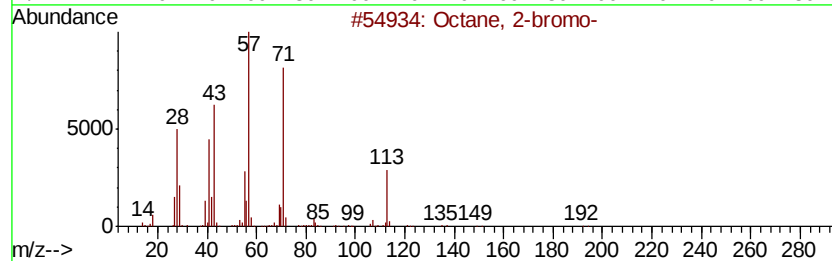
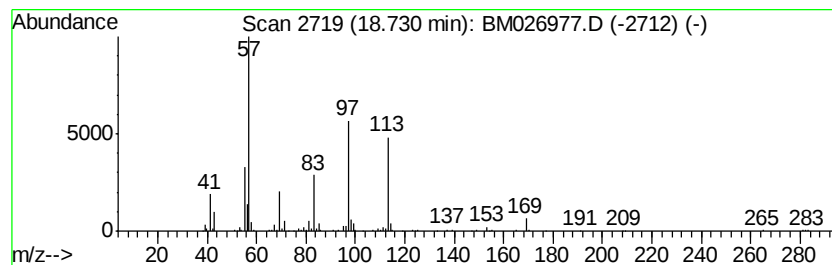
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.73	5.96 ng/ul	268366	Phenanthrene-d10	17.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octane, 2-bromo-	192 C8H17Br	000557-35-7	38
2	Sulfurous acid, cyclohexylmethyl...	276 C14H28O3S	1000309-21-7	38
3	2,4,4-Trimethyl-1-pentanol, trif...	226 C10H17F3O2	1000365-19-5	38
4	2,4,4-Trimethyl-1-pentanol, pent...	276 C11H17F5O2	1000365-51-0	35
5	Sulfurous acid, cyclohexylmethyl...	234 C11H22O3S	1000309-21-3	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
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Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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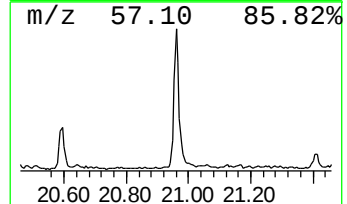
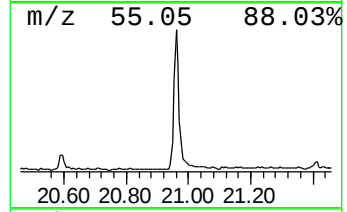
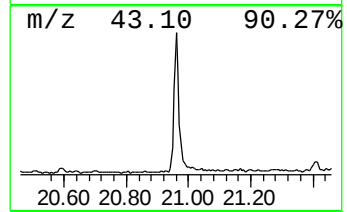
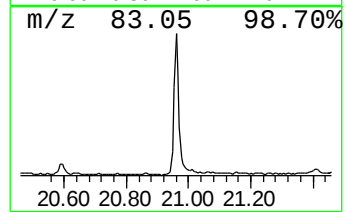
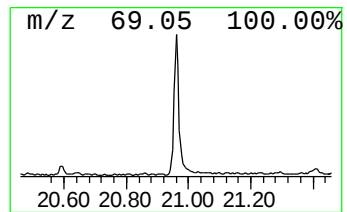
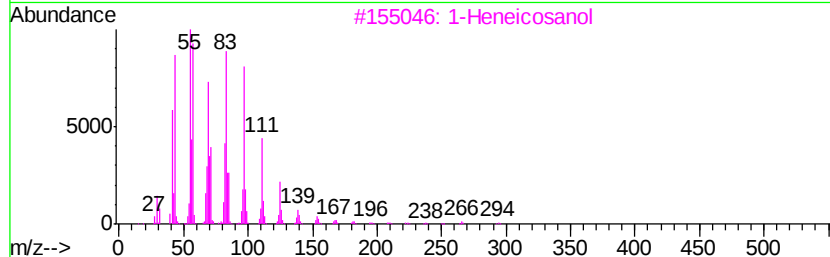
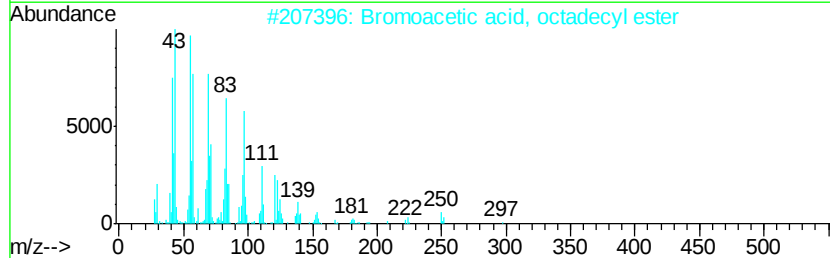
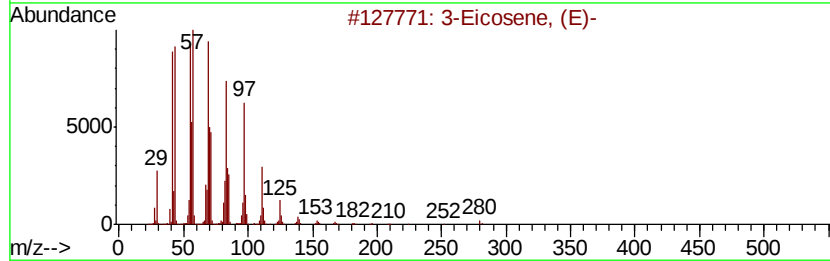
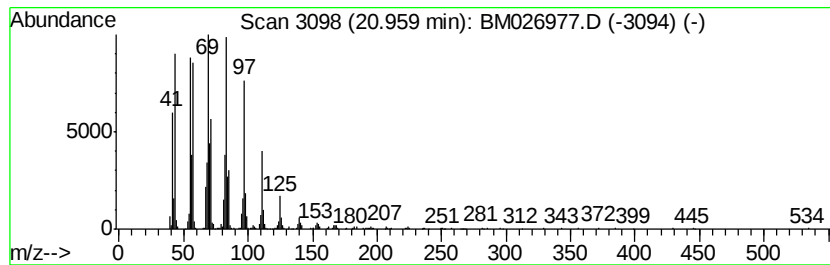
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	4.83 ng/ul	510048	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
2		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
3		1-Heneicosanol	312	C21H44O	015594-90-8	89
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	87
5		Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
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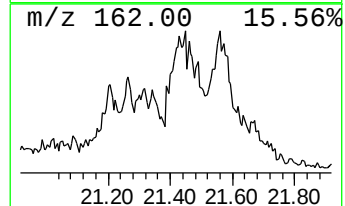
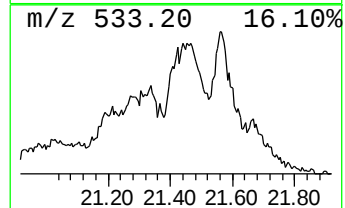
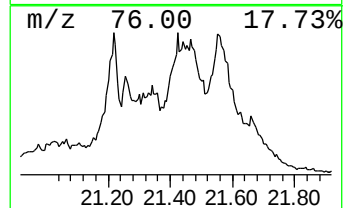
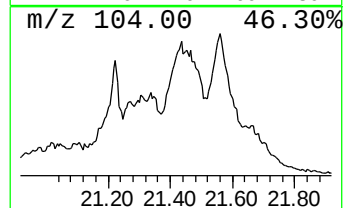
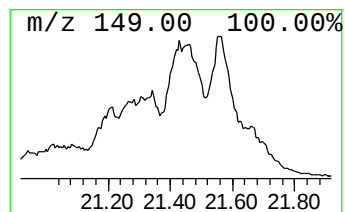
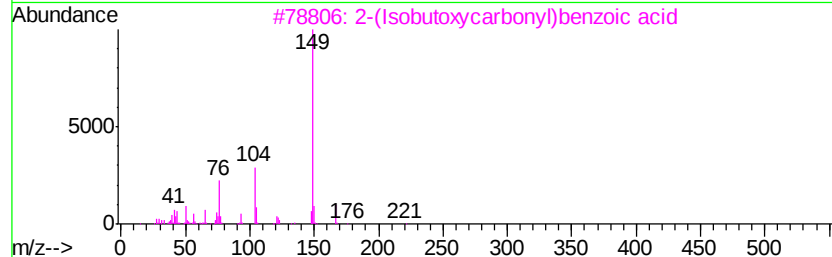
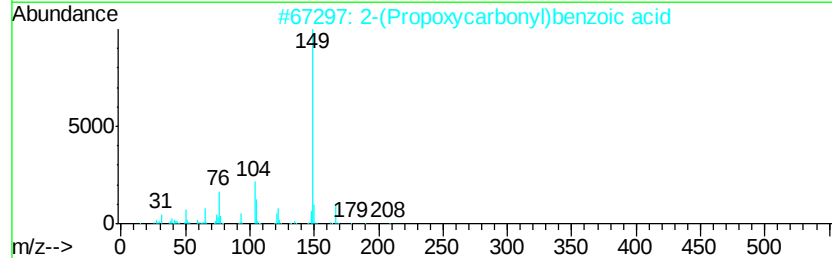
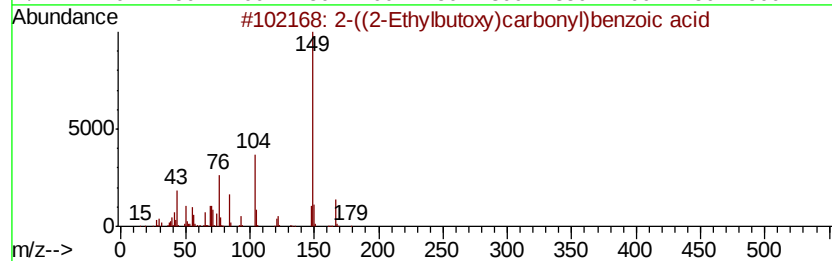
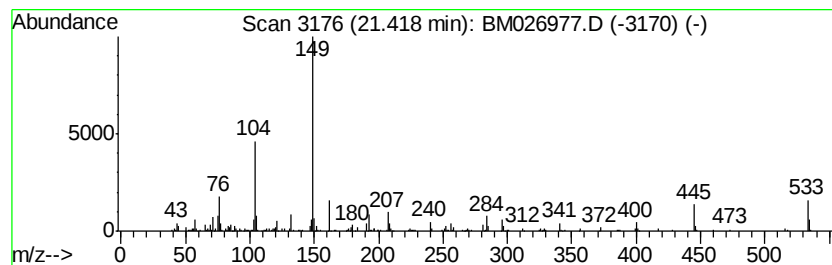
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-((2-Ethylbutoxy)carbonyl)... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.42	2.37 ng/ul	249684	Chrysene-d12	21.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-((2-Ethylbutoxy)carbonyl)benzo...	250	C14H18O4	1000373-62-8	53	
2	2-(Propoxycarbonyl)benzoic acid	208	C11H12O4	1000373-63-1	50	
3	2-(Isobutoxycarbonyl)benzoic acid	222	C12H14O4	1000373-52-8	45	
4	2-((Prop-2-ynyl)oxy)carbonyl)benz...	204	C11H8O4	1000373-53-6	45	
5	Phthalic acid, allyl ethyl ester	234	C13H14O4	033672-94-5	38	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
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Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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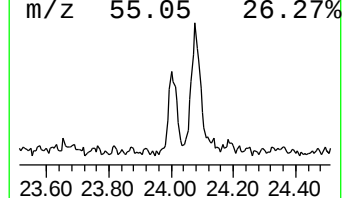
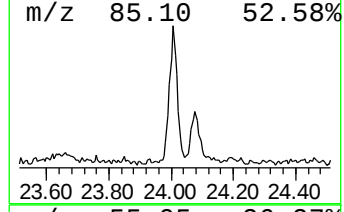
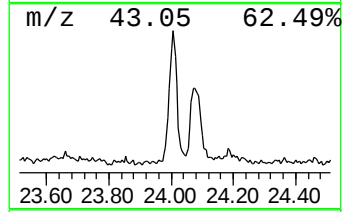
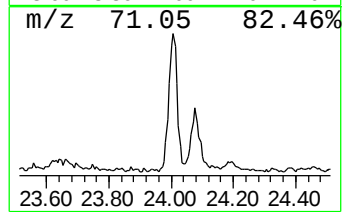
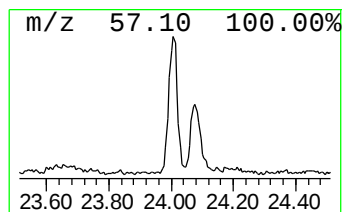
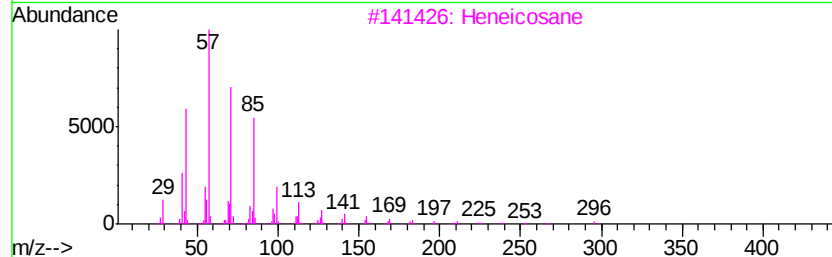
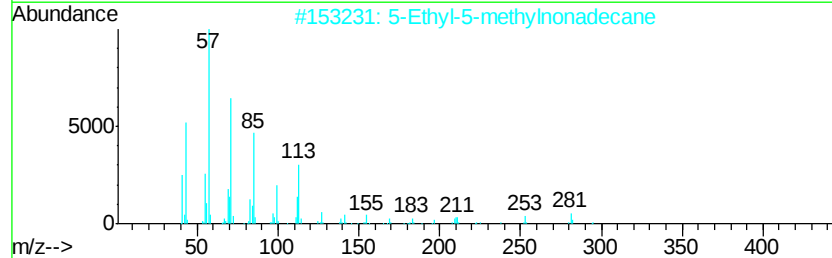
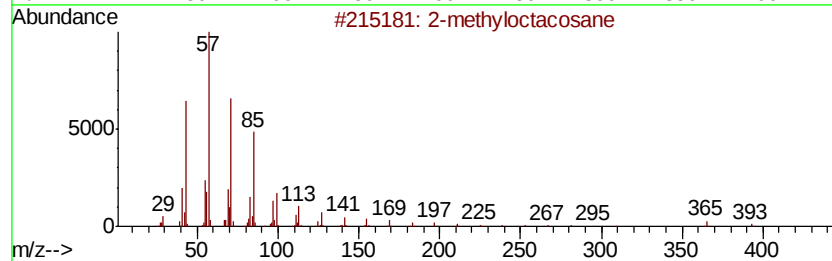
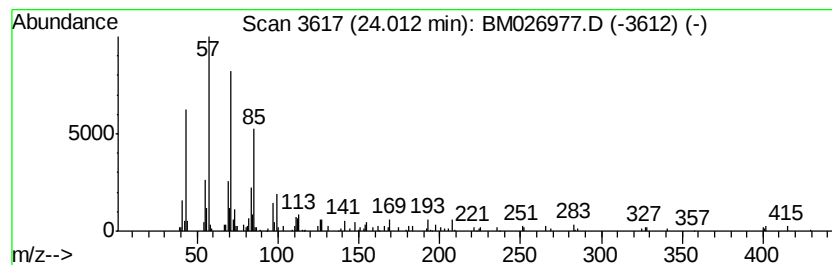
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.01	2.51 ng/ul	137822	Perylene-d12	23.43

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	87
2			5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	86
3			Heneicosane	296	C21H44	000629-94-7	80
4			Hexacosane	366	C26H54	000630-01-3	74
5			Tetracosane	338	C24H50	000646-31-1	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026977.D
Acq On : 03 Aug 2020 12:44
Operator : JU/CG
Sample : L3424-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0S60

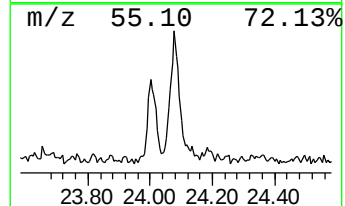
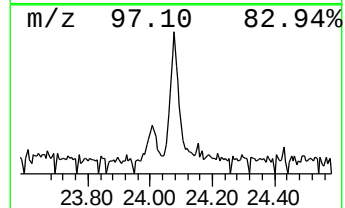
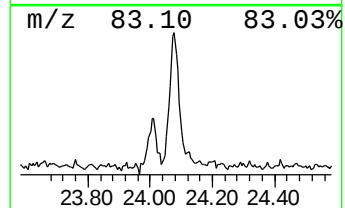
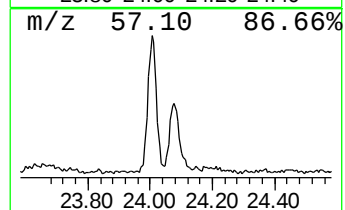
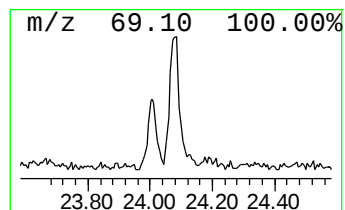
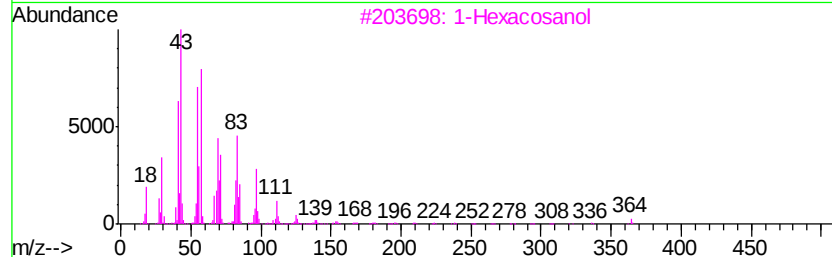
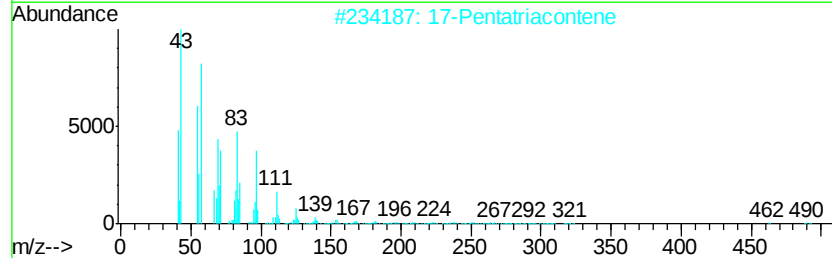
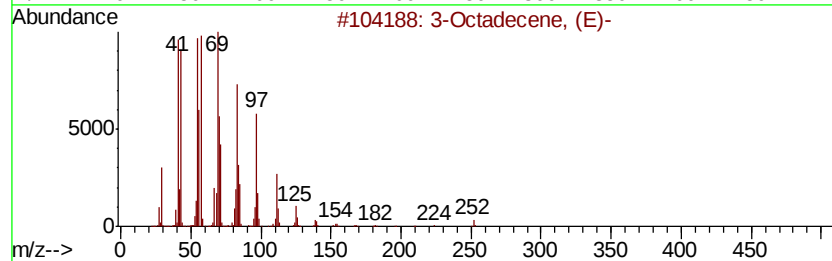
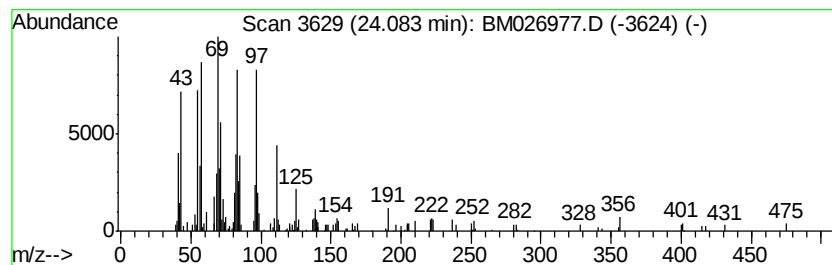
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	2.68 ng/ul	147246	Perylene-d12	23.43

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octadecene, (E)-	252	C18H36	007206-19-1	90
2			17-Pentatriacontene	491	C35H70	006971-40-0	74
3			1-Hexacosanol	382	C26H54O	000506-52-5	72
4			Hexadecane, 1-(ethenvloxy)-	268	C18H36O	000822-28-6	53
5			1-Heneicosanol	312	C21H44O	015594-90-8	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080320\
 Data File : BM026977.D
 Acq On : 03 Aug 2020 12:44
 Operator : JU/CG
 Sample : L3424-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0S60

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.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
(DEL) Alkane: Cyc...	2.90	9.4	ng/ul	191429	1	7.66	407223	20.0
Butane, 2-methoxy...	2.97	16.3	ng/ul	331700	1	7.66	407223	20.0
2-Pentanone, 4-hy...	4.83	10.3	ng/ul	210616	1	7.66	407223	20.0
Naphthalene, 1-me...	12.32	25.2	ng/ul	721142	2	10.44	572896	20.0
1,1'-Biphenyl, 4-...	17.51	22.7	ng/ul	1022390	4	17.04	901050	20.0
n-Hexadecanoic acid	17.95	3.4	ng/ul	154189	4	17.04	901050	20.0
unknown-01	18.73	6.0	ng/ul	268366	4	17.04	901050	20.0
(DEL) Alkane: Str...	20.96	4.8	ng/ul	510048	5	21.22	2110480	20.0
2-((2-Ethylbutoxy...	21.42	2.4	ng/ul	249684	5	21.22	2110480	20.0
(DEL) Alkane: Str...	24.01	2.5	ng/ul	137822	6	23.43	1098450	20.0
(DEL) Alkane: Str...	24.08	2.7	ng/ul	147246	6	23.43	1098450	20.0