

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
 Data File : BM021633.D
 Acq On : 25 Jul 2019 11:16
 Operator : HP/JU
 Sample : K3764-10
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A52S7

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-BM072319MA.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	4.269	214	218	228	rBV	519890	758523	9.38%	0.649%
2	4.863	315	319	329	rVB	378772	549688	6.80%	0.470%
3	6.875	656	661	675	rVB	407472	704479	8.71%	0.603%
4	7.051	686	691	699	rVB	288852	469988	5.81%	0.402%
5	7.140	701	706	716	rVV	424102	647664	8.01%	0.554%
6	7.234	717	722	725	rVV	416791	707432	8.75%	0.606%
7	7.269	725	728	737	rVB	484490	712653	8.81%	0.610%
8	7.698	796	801	809	rVB	329303	545393	6.74%	0.467%
9	8.234	884	892	897	rBV2	266574	670547	8.29%	0.574%
10	8.281	897	900	908	rVB2	31166	47395	0.59%	0.041%
11	8.410	916	922	932	rBV	497792	814004	10.06%	0.697%
12	8.510	933	939	953	rVB	300268	501406	6.20%	0.429%
13	8.763	976	982	989	rVB	413047	674050	8.33%	0.577%
14	8.863	993	999	1003	rVV	424071	726042	8.98%	0.621%
15	8.910	1003	1007	1016	rVB	457192	747744	9.25%	0.640%
16	9.581	1115	1121	1131	rVB	323231	530916	6.56%	0.454%
17	9.910	1171	1177	1190	rVB	434060	676866	8.37%	0.579%
18	10.110	1204	1211	1213	rBV	641005	1072577	13.26%	0.918%
19	10.480	1268	1274	1278	rVV	534747	879815	10.88%	0.753%
20	10.533	1278	1283	1293	rVB	676460	1142272	14.12%	0.978%
21	10.639	1294	1301	1314	rBV	481375	873413	10.80%	0.748%
22	10.798	1321	1328	1335	rVB	478601	764389	9.45%	0.654%
23	12.027	1531	1537	1542	rBV	20949	37823	0.47%	0.032%
24	12.075	1543	1545	1550	rVB3	7941	9127	0.11%	0.008%
25	12.280	1574	1580	1588	rBV2	28057	50053	0.62%	0.043%
26	12.369	1588	1595	1608	rVV	960956	1511277	18.69%	1.294%
27	12.480	1608	1614	1617	rVV	346546	517874	6.40%	0.443%
28	12.516	1617	1620	1630	rVB	409456	625048	7.73%	0.535%
29	12.769	1657	1663	1670	rBV	1062796	1500169	18.55%	1.284%
30	12.845	1670	1676	1691	rVB	640384	1100852	13.61%	0.942%
31	13.145	1723	1727	1731	rBV4	8840	12581	0.16%	0.011%
32	13.216	1732	1739	1749	rVB	1012415	1526083	18.87%	1.306%
33	13.763	1826	1832	1836	rVV	1242933	1688768	20.88%	1.445%
34	13.810	1836	1840	1851	rVB	1504204	2011103	24.87%	1.721%

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35	13.945	1856	1863	1869	rBV	353814	489939	6.06%	0.419%
36	14.039	1872	1879	1882	rVV	1449433	2349807	29.05%	2.011%
37	14.069	1882	1884	1892	rVB	948683	1075365	13.30%	0.920%
38	14.345	1924	1931	1937	rBV2	949661	1416043	17.51%	1.212%
39	14.410	1937	1942	1949	rVB	1023709	1459707	18.05%	1.249%
40	14.574	1963	1970	1975	rBV	359775	621391	7.68%	0.532%
41	14.627	1975	1979	1991	rVV	532525	785324	9.71%	0.672%
42	14.745	1992	1999	2008	rVB2	1324414	2017232	24.94%	1.727%
43	15.174	2066	2072	2081	rBV	1099767	1438990	17.79%	1.232%
44	15.345	2093	2101	2105	rBV	1939772	2722045	33.66%	2.330%
45	15.398	2105	2110	2119	rVV2	1967155	2732820	33.79%	2.339%
46	15.480	2119	2124	2137	rVV	383678	555393	6.87%	0.475%
47	15.580	2137	2141	2146	rVB	20242	26305	0.33%	0.023%
48	16.292	2256	2262	2268	rBV	909717	1193881	14.76%	1.022%
49	16.392	2273	2279	2285	rVB	707888	998081	12.34%	0.854%
50	16.751	2334	2340	2353	rBV	766944	1160868	14.35%	0.994%
51	16.845	2353	2356	2359	rVV3	8461	10092	0.12%	0.009%
52	16.898	2359	2365	2368	rVB3	6170	10846	0.13%	0.009%
53	17.098	2393	2399	2405	rBV2	1398948	1880949	23.26%	1.610%
54	17.198	2410	2416	2419	rVV	2247919	3219960	39.81%	2.756%
55	17.233	2419	2422	2434	rVB	2511549	3277222	40.52%	2.805%
56	17.509	2460	2469	2479	rBV	1961474	2824375	34.92%	2.417%
57	17.709	2499	2503	2506	rBV2	20015	22694	0.28%	0.019%
58	17.992	2546	2551	2558	rBV	233241	352075	4.35%	0.301%
59	18.068	2558	2564	2574	rVB	1885167	2267292	28.03%	1.941%
60	18.280	2596	2600	2604	rVB3	10589	13984	0.17%	0.012%
61	18.533	2638	2643	2652	rBV2	62288	102237	1.26%	0.088%
62	18.851	2695	2697	2705	rBV8	9683	21734	0.27%	0.019%
63	18.915	2705	2708	2711	rBV3	14161	18192	0.22%	0.016%
64	19.133	2741	2745	2749	rBV	41905	62874	0.78%	0.054%
65	19.274	2765	2769	2777	rBV	161584	245624	3.04%	0.210%
66	19.386	2785	2788	2791	rVB	19780	23143	0.29%	0.020%
67	19.498	2801	2807	2822	rVB	3144506	4194876	51.87%	3.590%
68	20.033	2895	2898	2903	rVB4	26595	33025	0.41%	0.028%
69	20.433	2961	2966	2976	rBV	2245612	2443176	30.21%	2.091%
70	20.533	2980	2983	2988	rVB	53021	67835	0.84%	0.058%
71	20.997	3058	3062	3075	rBV	373790	654495	8.09%	0.560%

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72	21.203	3092	3097	3104	rBV	1575578	1847389	22.84%	1.581%
73	21.280	3104	3110	3122	rVV2	3542585	6776155	83.78%	5.800%
74	21.433	3132	3136	3143	rBV	523246	571013	7.06%	0.489%
75	21.868	3205	3210	3219	rBV	988540	1260420	15.58%	1.079%
76	22.086	3242	3247	3257	rVB	3307342	3873439	47.89%	3.315%
77	22.333	3284	3289	3297	rBV	1564363	2090670	25.85%	1.789%
78	22.839	3369	3375	3377	rBV	1871743	2813129	34.78%	2.408%
79	22.862	3377	3379	3383	rVV	1348801	2055816	25.42%	1.760%
80	22.915	3383	3388	3402	rVB	3624918	6062693	74.96%	5.189%
81	23.397	3463	3470	3485	rBV	4513024	8087918	100.00%	6.923%
82	23.544	3488	3495	3506	rVB	1519801	2960708	36.61%	2.534%
83	24.038	3573	3579	3589	rBV	1427221	2652576	32.80%	2.270%
84	24.780	3699	3705	3714	rVB	842163	1852004	22.90%	1.585%
85	25.644	3846	3852	3864	rVB	416746	1049168	12.97%	0.898%
86	26.497	3987	3997	4013	rVB	1119564	3626563	44.84%	3.104%
87	26.662	4020	4025	4036	rVB2	250896	656226	8.11%	0.562%

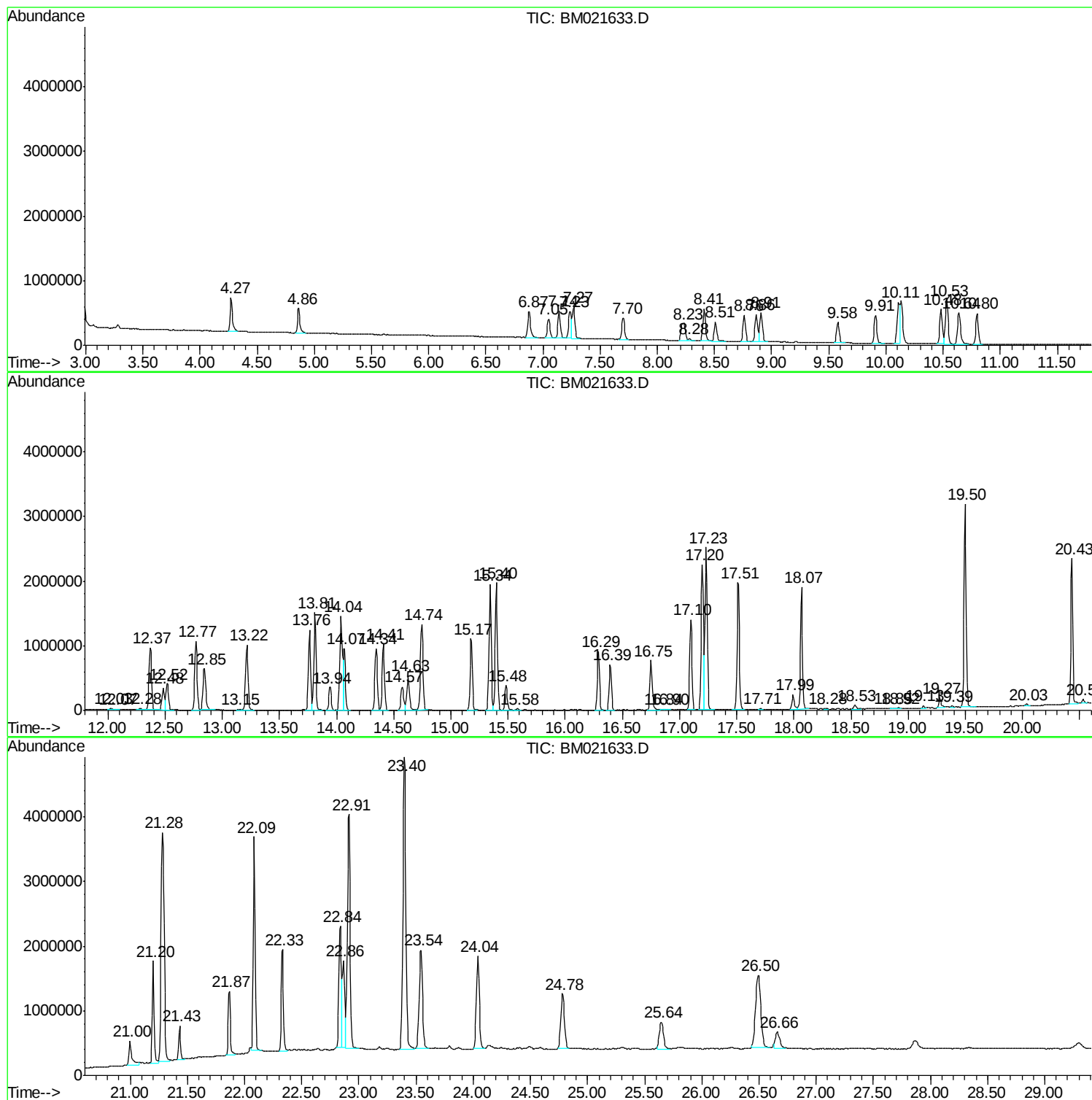
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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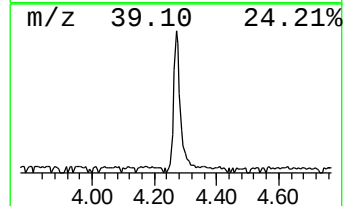
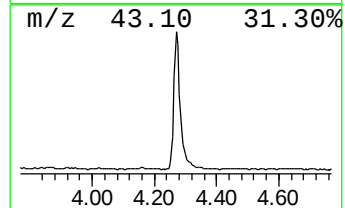
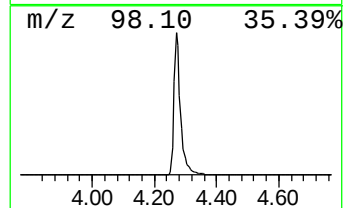
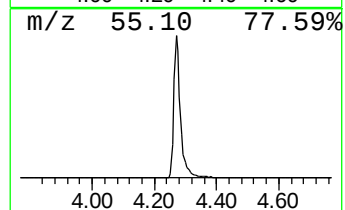
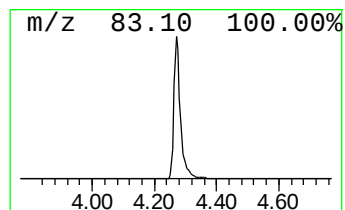
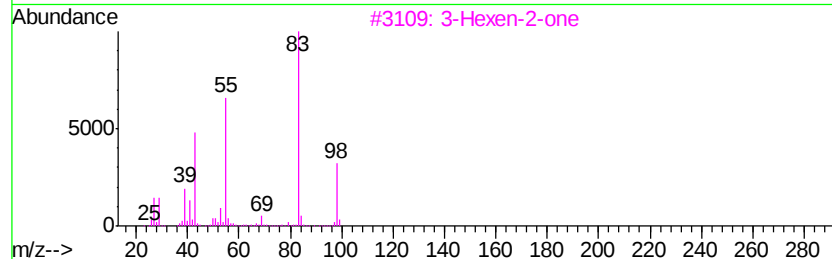
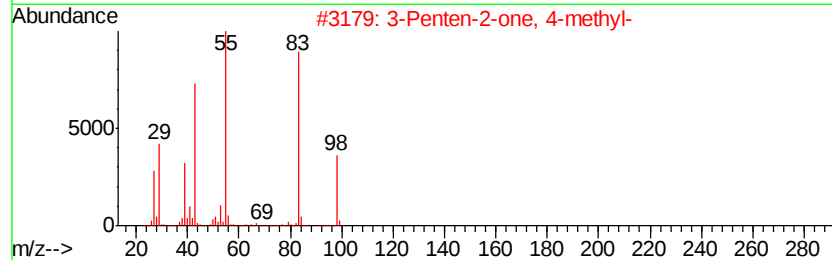
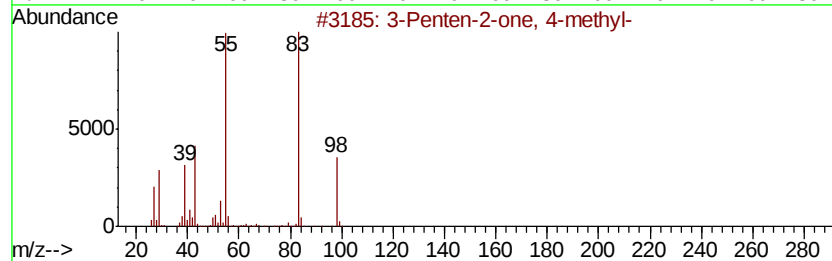
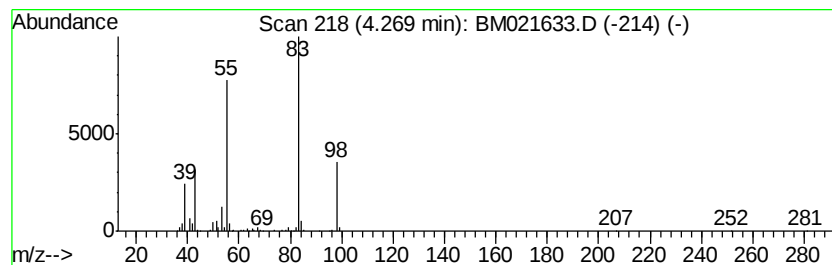
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.27	27.82 ng/ul	758523	1,4-Dichlorobenzene-d4	7.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3			3-Hexen-2-one	98	C6H10O	000763-93-9	91
4			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
5			3-Hexen-2-one	98	C6H10O	000763-93-9	90



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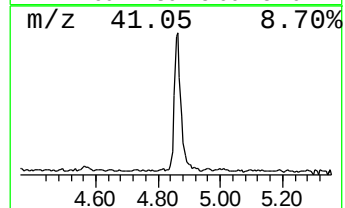
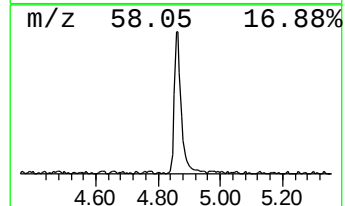
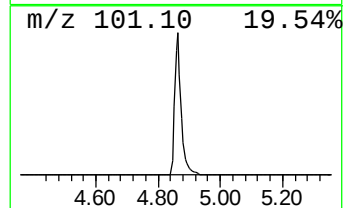
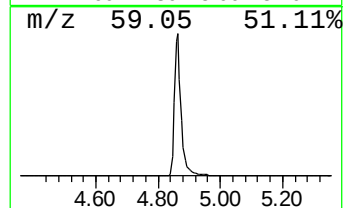
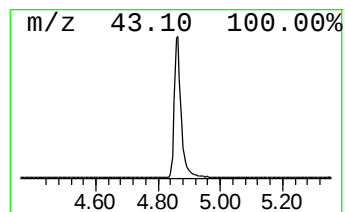
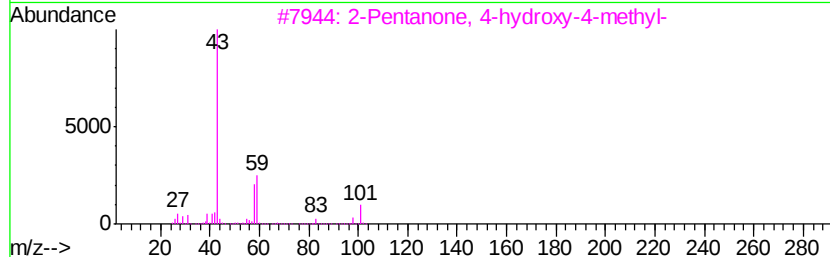
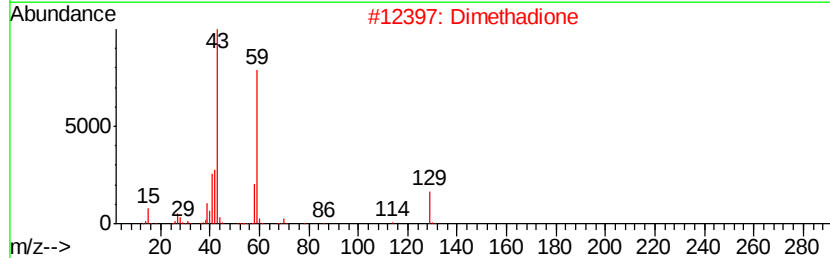
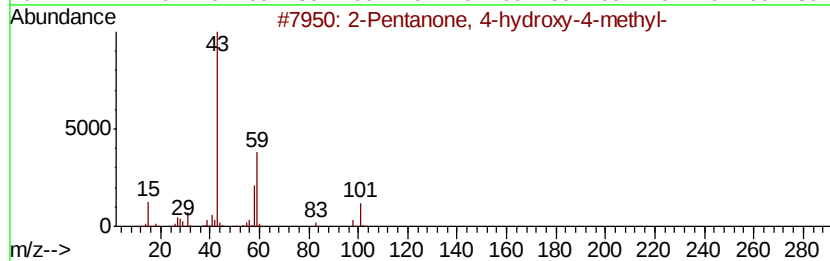
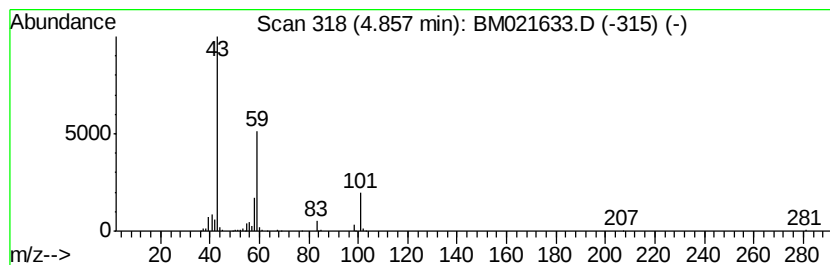
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	20.16 ng/ul	549688	1,4-Dichlorobenzene-d4	7.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Dimethadione	129	C5H7NO3	000695-53-4	40
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
5		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28



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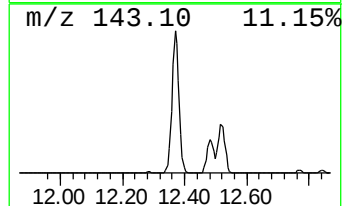
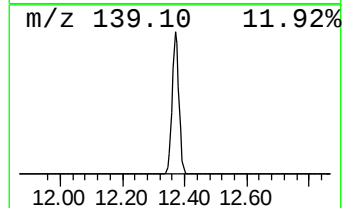
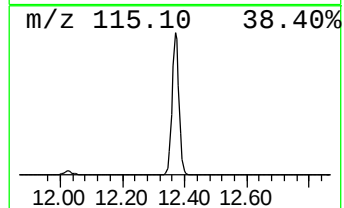
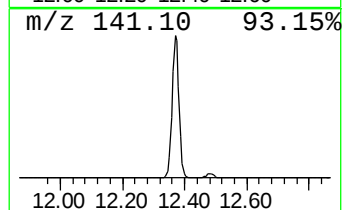
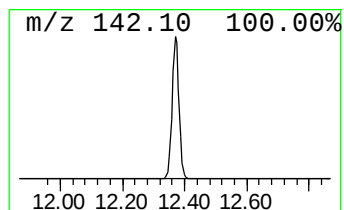
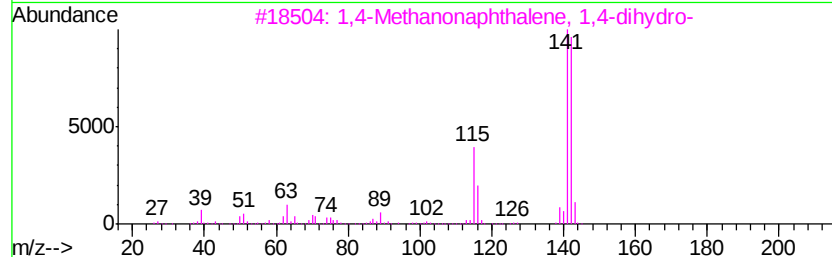
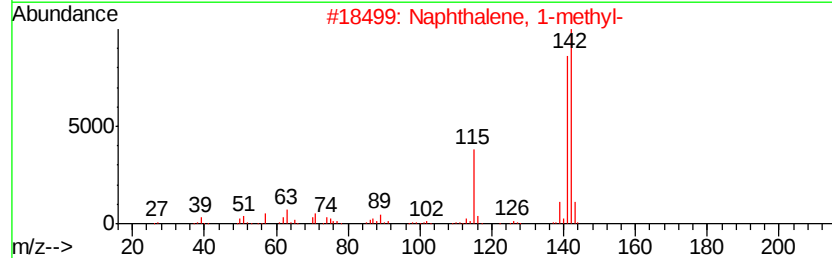
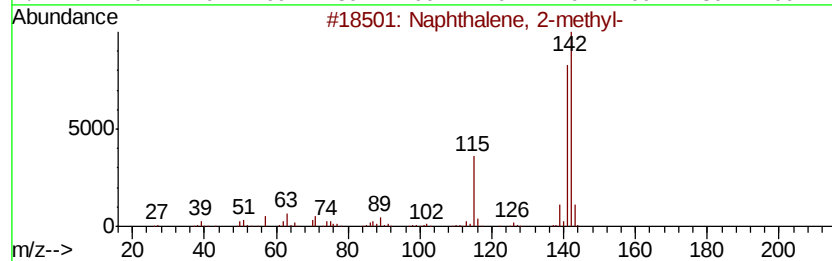
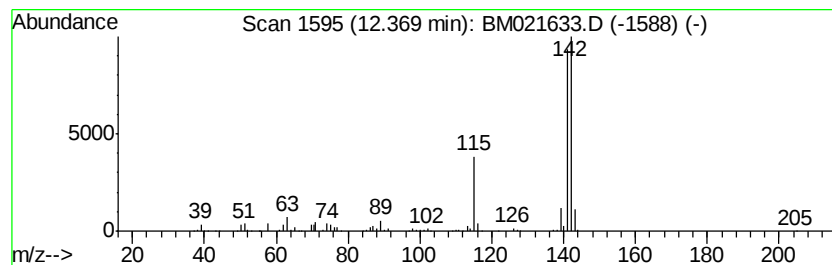
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	34.35 ng/ul	1511280	Naphthalene-d8	10.48

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



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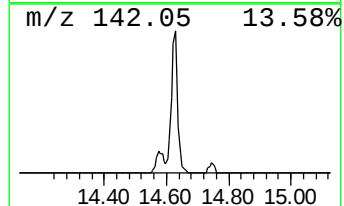
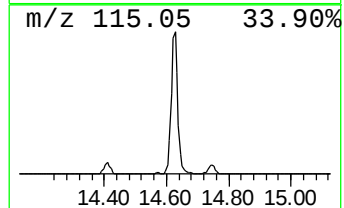
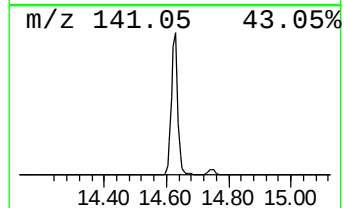
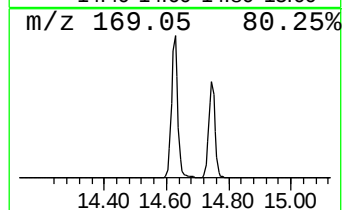
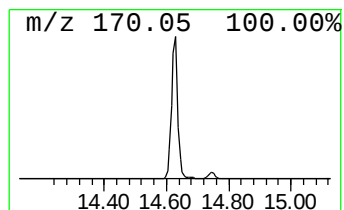
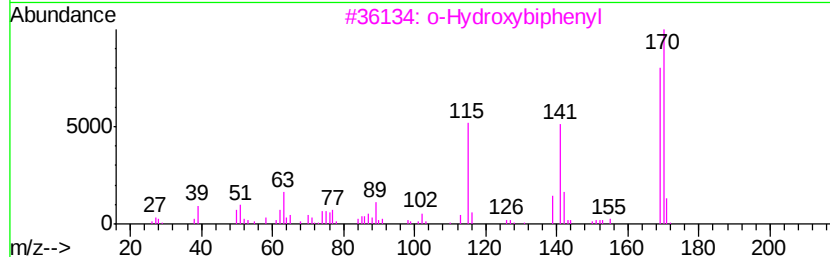
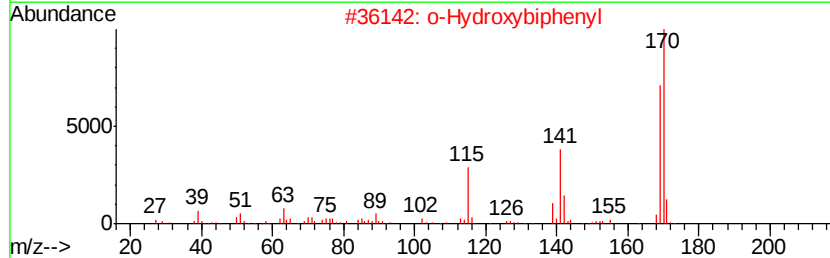
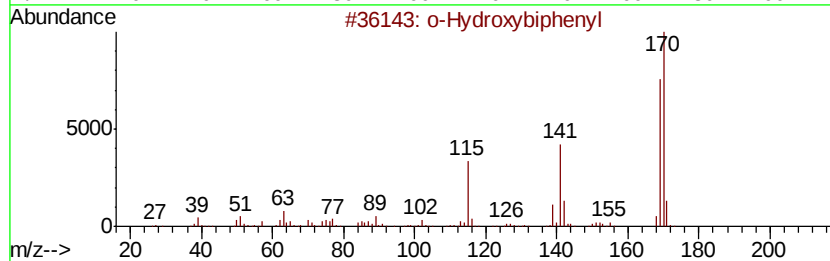
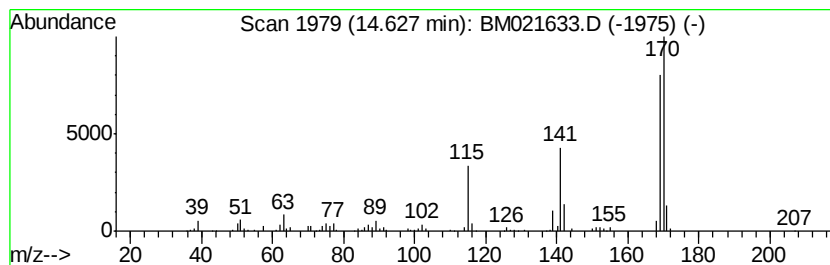
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 o-Hydroxybiphenyl Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	11.09 ng/ul	785324	Acenaphthene-d10	14.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Hydroxybiphenyl	170 C12H10O	000090-43-7 97
2		o-Hydroxybiphenyl	170 C12H10O	000090-43-7 96
3		o-Hydroxybiphenyl	170 C12H10O	000090-43-7 96
4		o-Hydroxybiphenyl	170 C12H10O	000090-43-7 87
5		[1,1'-Biphenyl]-3-ol	170 C12H10O	000580-51-8 58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
Operator : HP/JU
Sample : K3764-10
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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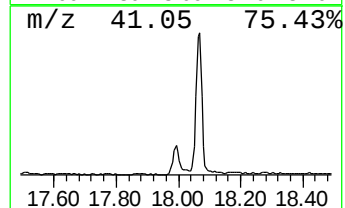
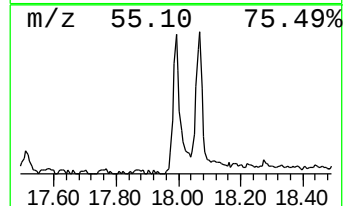
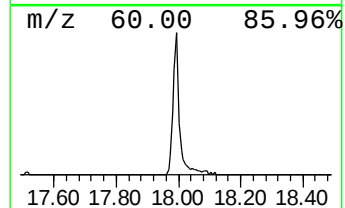
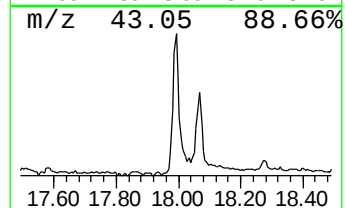
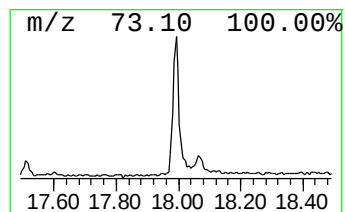
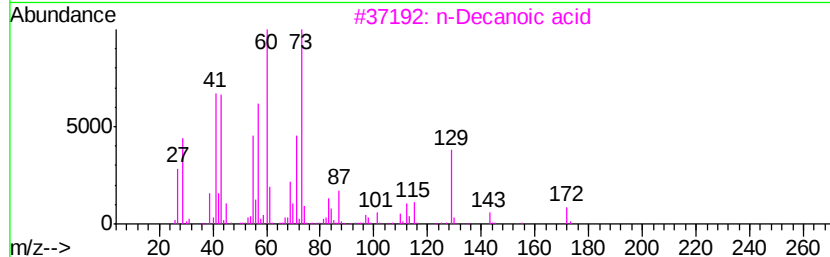
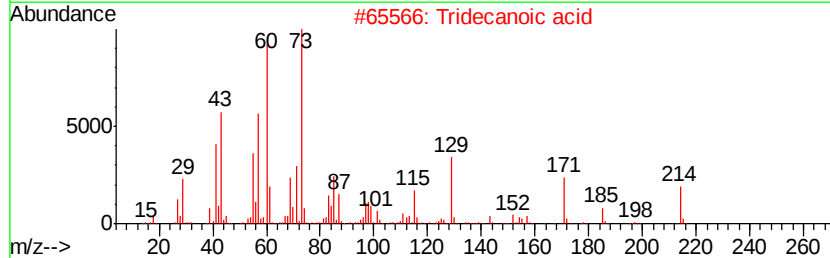
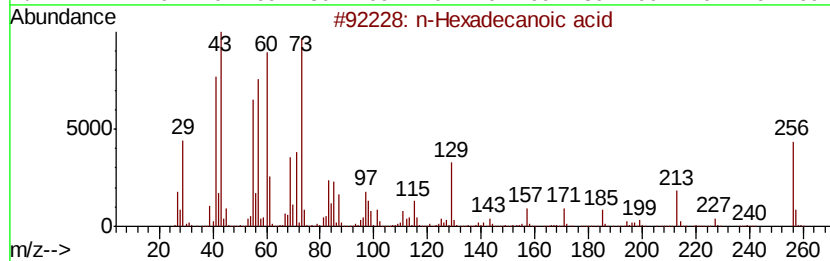
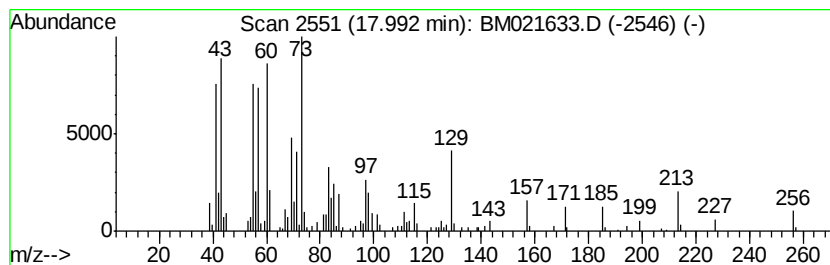
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.99	3.74 ng/ul	352075	Phenanthrene-d10	17.10

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	81
3		n-Decanoic acid	172	C10H20O2	000334-48-5	76
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
Operator : HP/JU
Sample : K3764-10
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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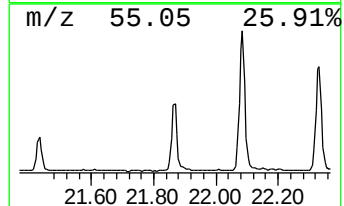
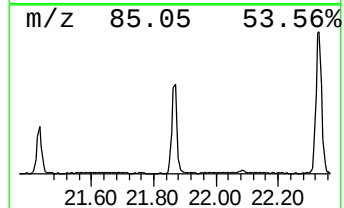
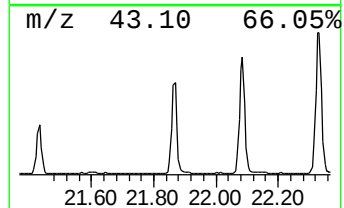
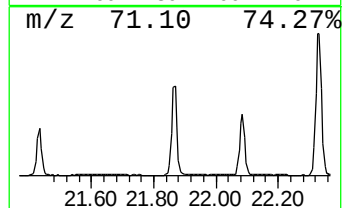
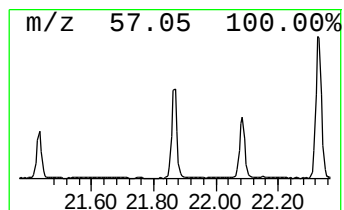
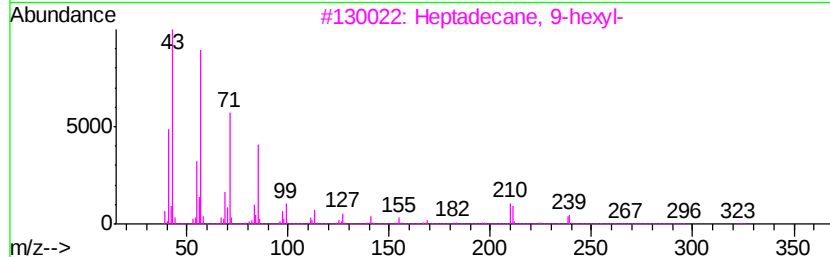
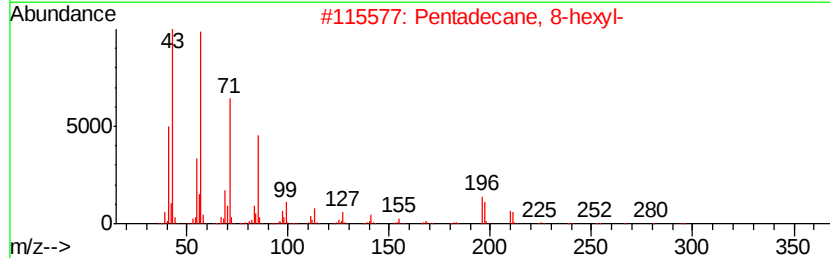
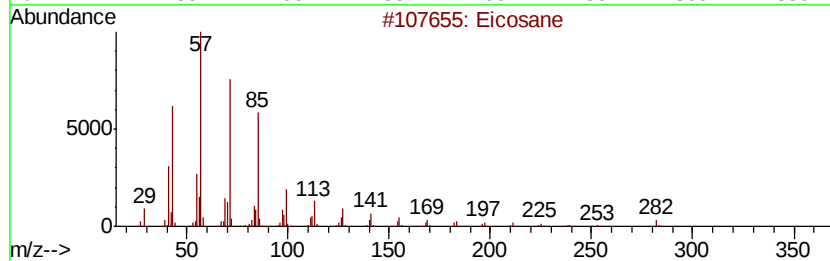
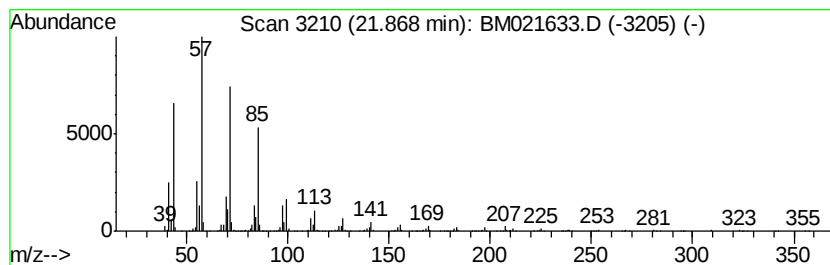
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.87	3.72 ng/ul	1260420	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Hexadecane	226	C16H34	000544-76-3	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
Operator : HP/JU
Sample : K3764-10
Misc :
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Instrument :
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ClientSampleId :
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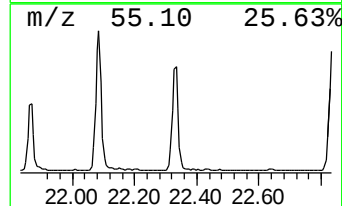
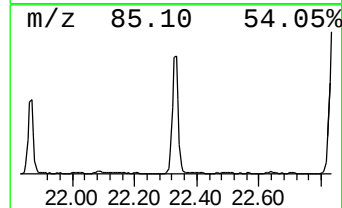
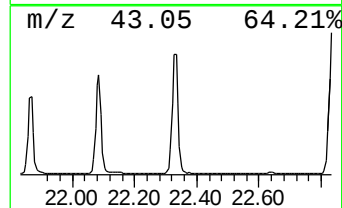
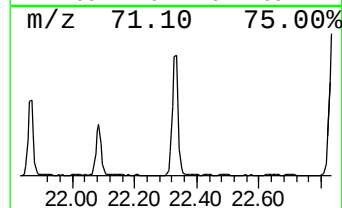
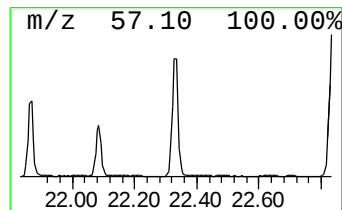
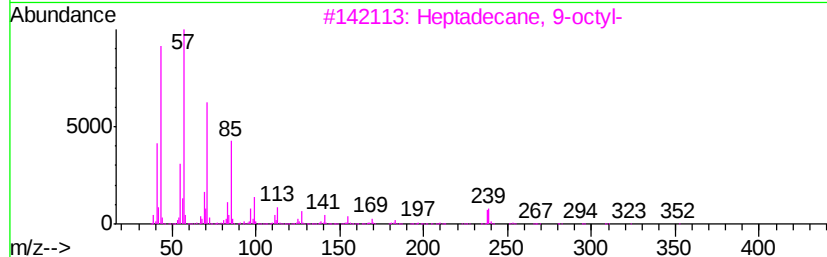
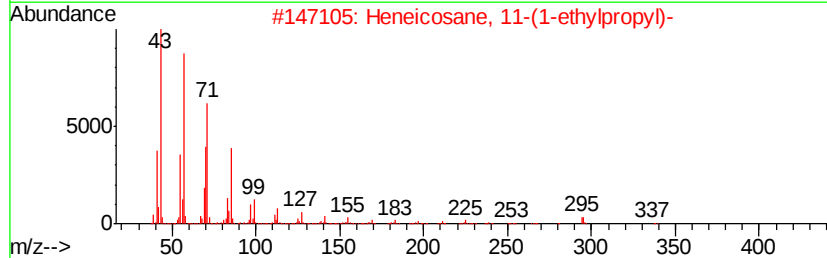
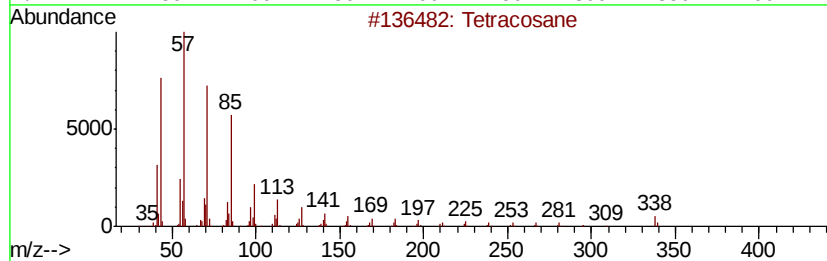
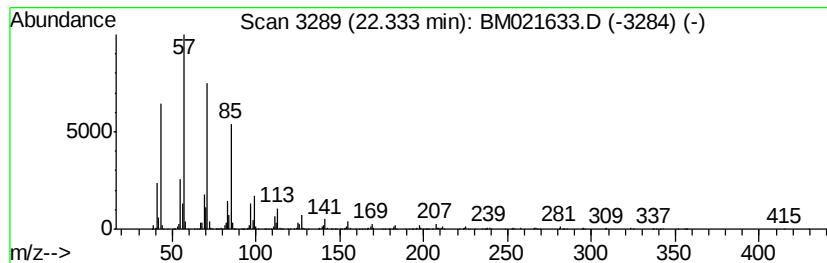
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.33	6.17 ng/ul	2090670	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
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Misc :
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Instrument :
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ClientSampleId :
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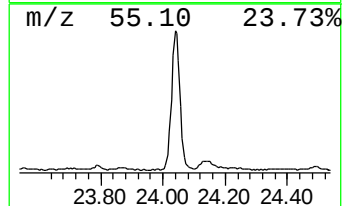
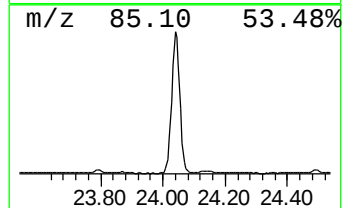
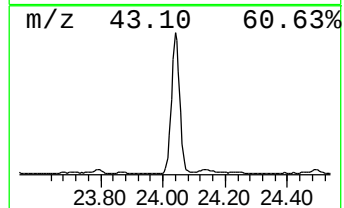
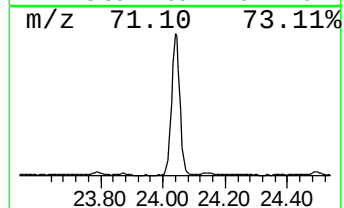
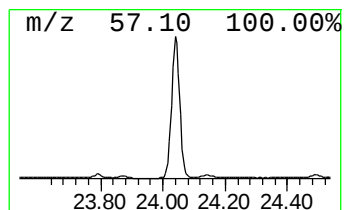
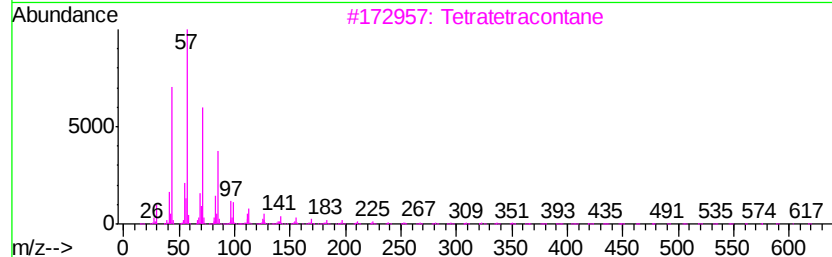
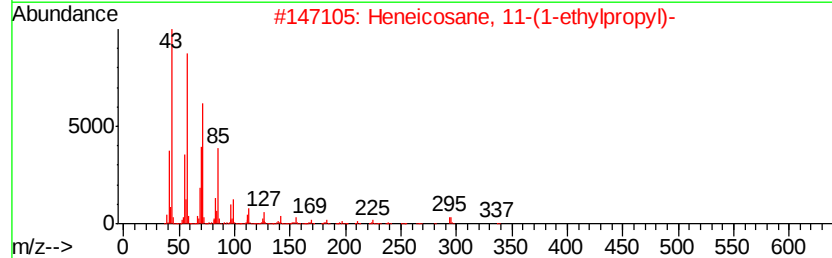
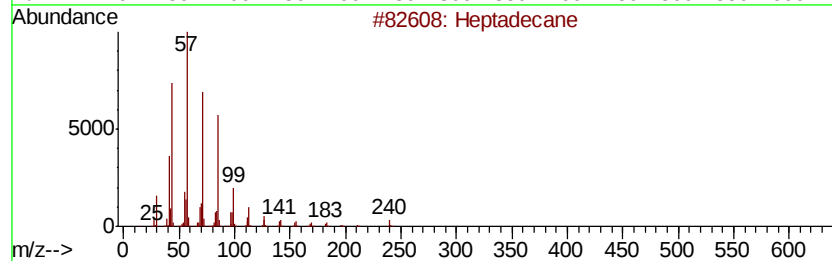
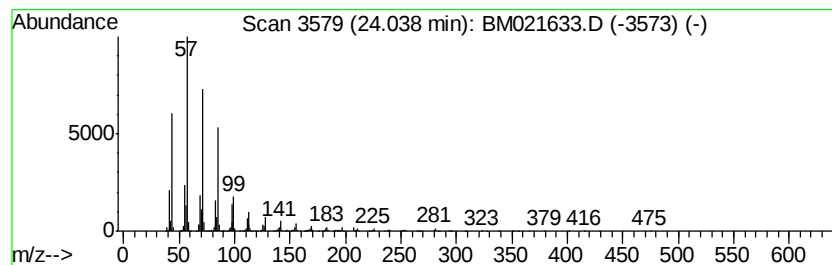
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.04	17.92 ng/ul	2652580	Perylene-d12	23.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
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Instrument :
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ClientSampleId :
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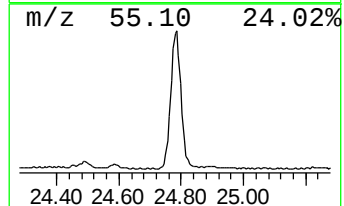
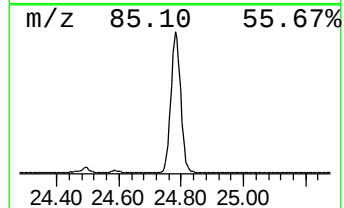
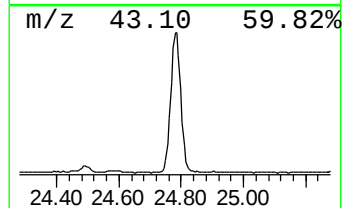
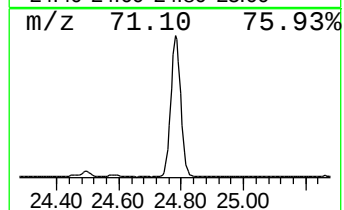
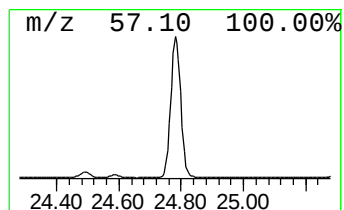
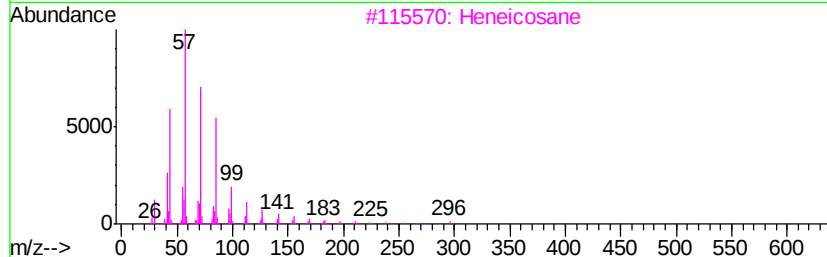
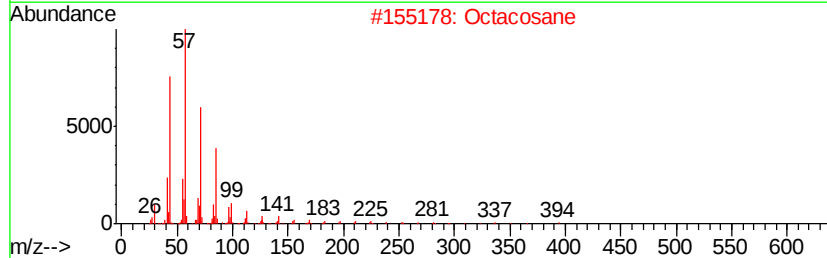
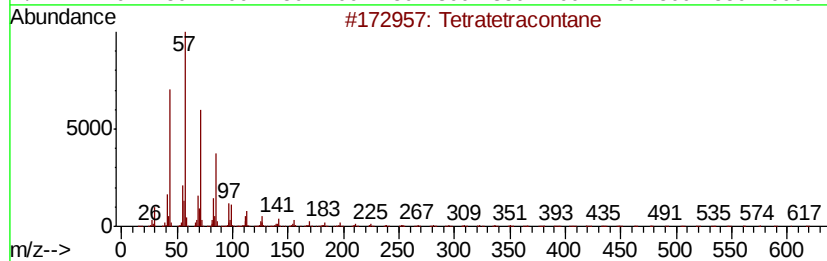
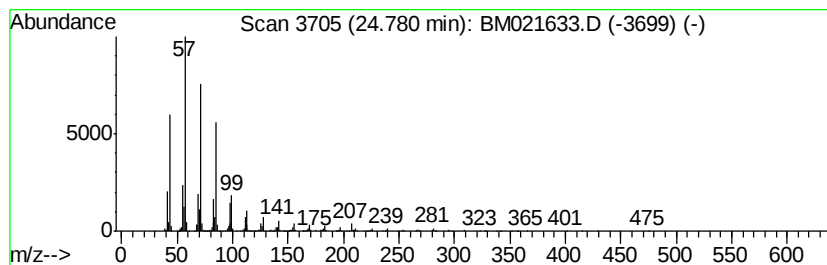
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.78	12.51 ng/ul	1852000	Perylene-d12	23.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptacosane	380	C27H56	000593-49-7	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
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ClientSampleId :
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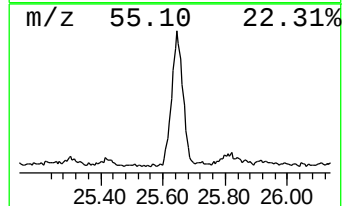
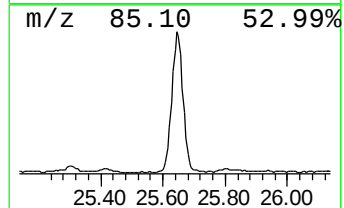
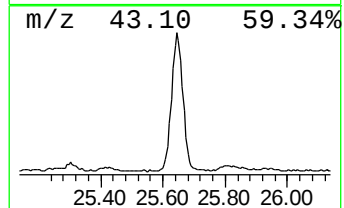
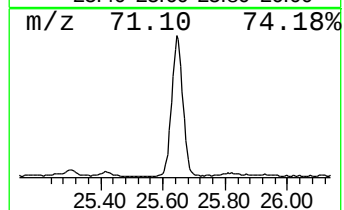
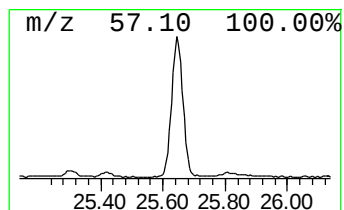
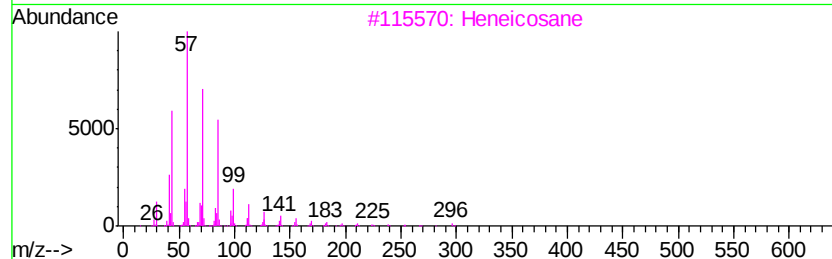
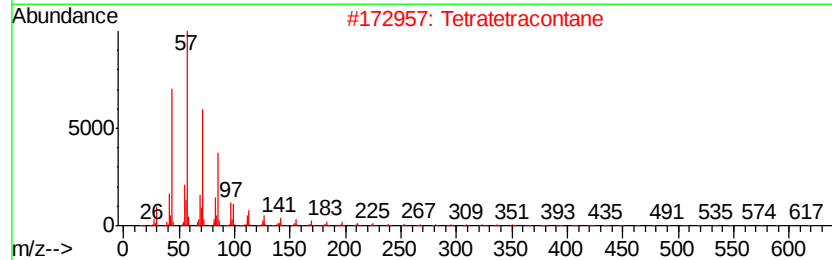
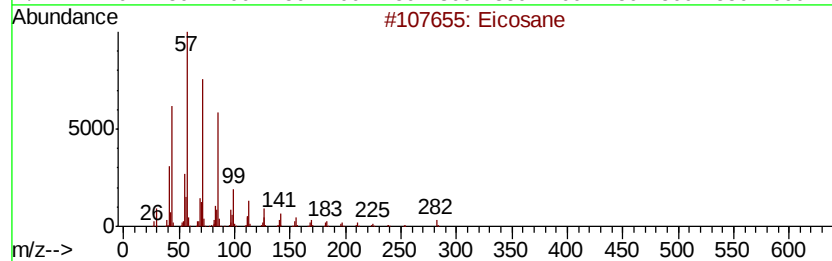
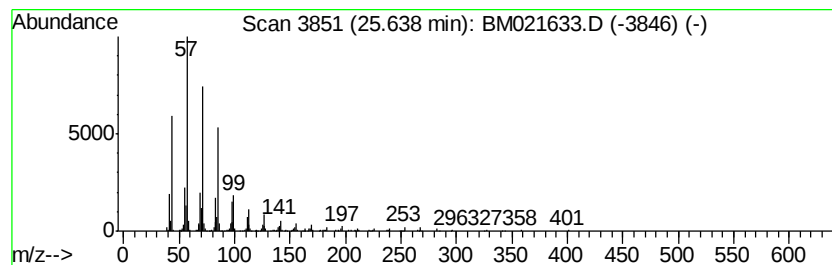
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.64	7.09 ng/ul	1049170	Perylene-d12	23.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
Operator : HP/JU
Sample : K3764-10
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A52S7

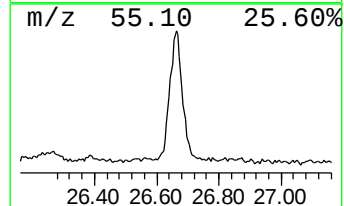
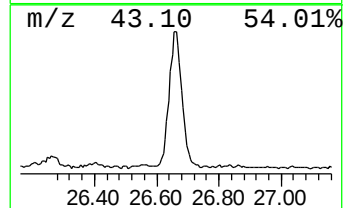
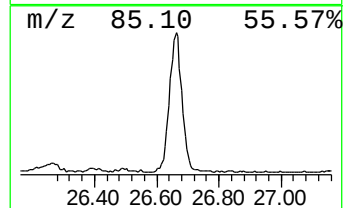
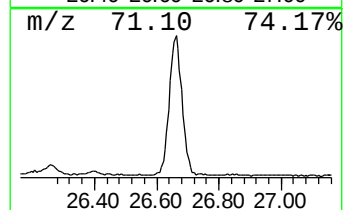
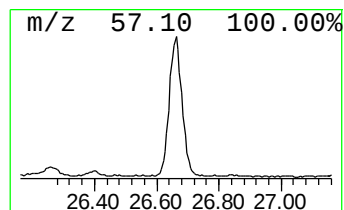
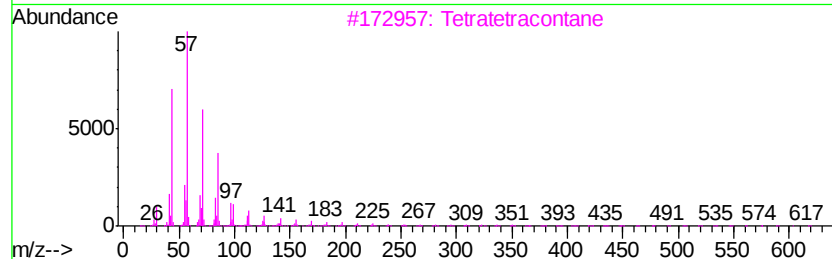
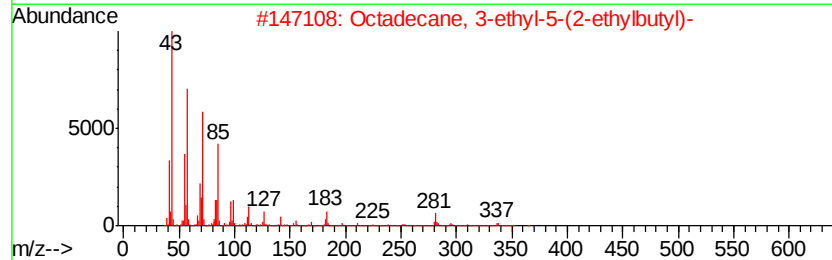
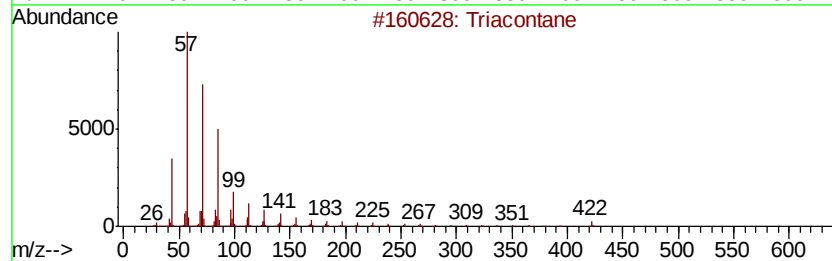
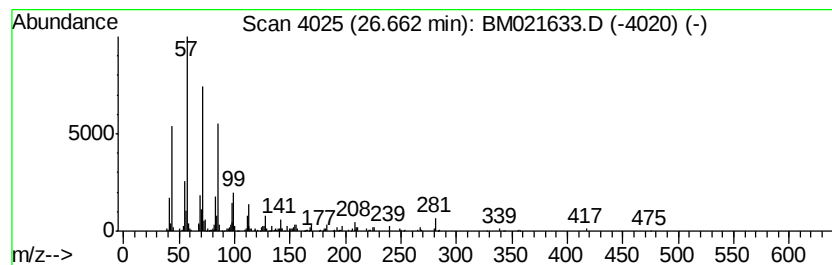
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.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.
26.66	4.43 ng/ul	656226	Perylene-d12	23.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Triacontane	422	C30H62	000638-68-6	90
2		Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	90
3		Tetratetracontane	619	C44H90	007098-22-8	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM072519\
Data File : BM021633.D
Acq On : 25 Jul 2019 11:16
Operator : HP/JU
Sample : K3764-10
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A52S7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.27	27.8	ng/ul	758523	1	7.70	545393	20.0
2-Pentanone, 4-hy...	4.86	20.2	ng/ul	549688	1	7.70	545393	20.0
Naphthalene, 1-me...	12.37	34.4	ng/ul	1511280	2	10.48	879815	20.0
o-Hydroxybiphenyl	14.63	11.1	ng/ul	785324	3	14.35	1416040	20.0
n-Hexadecanoic acid	17.99	3.7	ng/ul	352075	4	17.10	1880950	20.0
(DEL) Alkane: Str...	21.87	3.7	ng/ul	1260420	5	21.29	6776160	20.0
(DEL) Alkane: Str...	22.33	6.2	ng/ul	2090670	5	21.29	6776160	20.0
(DEL) Alkane: Str...	24.04	17.9	ng/ul	2652580	6	23.54	2960710	20.0
(DEL) Alkane: Str...	24.78	12.5	ng/ul	1852000	6	23.54	2960710	20.0
(DEL) Alkane: Str...	25.64	7.1	ng/ul	1049170	6	23.54	2960710	20.0
(DEL) Alkane: Str...	26.66	4.4	ng/ul	656226	6	23.54	2960710	20.0