

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
 Data File : BM023320.D
 Acq On : 22 Oct 2019 12:29
 Operator : JU
 Sample : K5308-07
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X2J25

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-BM101419MA.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	3.169	28	31	38	rVB	75108	103137	0.29%	0.036%
2	3.893	150	154	158	rVB	31505	45668	0.13%	0.016%
3	4.216	204	209	218	rVB	480687	650069	1.83%	0.228%
4	4.804	304	309	319	rBV	475075	663767	1.87%	0.233%
5	6.798	640	648	651	rBV	3494281	5875574	16.51%	2.060%
6	6.993	674	681	690	rBV	921125	1433876	4.03%	0.503%
7	7.187	707	714	716	rBV	1341022	2052021	5.77%	0.719%
8	7.216	716	719	728	rVV	2810851	4377387	12.30%	1.534%
9	7.651	786	793	800	rBV	1235179	1949471	5.48%	0.683%
10	8.357	906	913	918	rBV	1487907	2306069	6.48%	0.808%
11	8.416	918	923	927	rVV	1673068	2590227	7.28%	0.908%
12	8.469	927	932	942	rVB	3021270	4749033	13.35%	1.665%
13	8.728	967	976	982	rBV	2180144	3594375	10.10%	1.260%
14	8.798	982	988	998	rVB	1107399	1820638	5.12%	0.638%
15	9.275	1064	1069	1076	rVB2	45553	71274	0.20%	0.025%
16	9.516	1103	1110	1119	rBV	870361	1415501	3.98%	0.496%
17	10.081	1193	1206	1216	rBV2	3865970	8378174	23.55%	2.937%
18	10.428	1258	1265	1269	rBV	1599401	2753213	7.74%	0.965%
19	10.481	1269	1274	1282	rVV	4729016	7711911	21.68%	2.703%
20	10.563	1282	1288	1305	rVB	1324048	2291201	6.44%	0.803%
21	11.345	1415	1421	1439	rBV	226912	559821	1.57%	0.196%
22	11.975	1520	1528	1533	rBV3	49616	98628	0.28%	0.035%
23	12.022	1533	1536	1542	rVB2	25940	38574	0.11%	0.014%
24	12.104	1542	1550	1559	rBV	5661324	8991430	25.27%	3.152%
25	12.322	1582	1587	1593	rVB	64891	100456	0.28%	0.035%
26	12.769	1658	1663	1668	rBV2	44502	68458	0.19%	0.024%
27	13.116	1717	1722	1724	rBV	69580	104909	0.29%	0.037%
28	13.163	1724	1730	1737	rVB	6047609	9119018	25.63%	3.196%
29	13.710	1816	1823	1828	rBV	2635279	3631832	10.21%	1.273%
30	13.757	1828	1831	1839	rVV	705262	939274	2.64%	0.329%
31	13.986	1862	1870	1876	rBV	2914700	4295703	12.07%	1.506%
32	14.292	1916	1922	1928	rBV	2372638	3612619	10.15%	1.266%
33	14.357	1928	1933	1941	rVB	4152300	6106732	17.16%	2.141%
34	14.492	1945	1956	1966	rBV	1227827	1810983	5.09%	0.635%

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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Title : SVOA CALIBRATION

35	14.727	1991	1996	2000	rVB7	27651	40275	0.11%	0.014%
36	15.180	2068	2073	2080	rVB3	33971	49401	0.14%	0.017%
37	15.292	2085	2092	2096	rBV	3921156	5507103	15.48%	1.930%
38	15.345	2096	2101	2107	rVV	4843712	6774750	19.04%	2.375%
39	15.421	2107	2114	2128	rVV3	3936256	6514598	18.31%	2.284%
40	15.533	2128	2133	2138	rVV2	53097	95939	0.27%	0.034%
41	16.457	2286	2290	2294	rBV2	61569	82151	0.23%	0.029%
42	16.521	2294	2301	2306	rBV	5848601	7723791	21.71%	2.707%
43	16.557	2306	2307	2311	rVB2	80296	71575	0.20%	0.025%
44	16.857	2351	2358	2364	rVB	114219	175635	0.49%	0.062%
45	17.039	2383	2389	2392	rBV2	3008683	4296469	12.08%	1.506%
46	17.080	2392	2396	2402	rVV	7124686	10724820	30.14%	3.759%
47	17.139	2402	2406	2415	rVB	4370321	6056423	17.02%	2.123%
48	17.251	2420	2425	2430	rVB8	31762	47887	0.13%	0.017%
49	17.445	2451	2458	2468	rBV	9451510	13343217	37.50%	4.677%
50	17.692	2494	2500	2505	rBV2	224282	316191	0.89%	0.111%
51	17.962	2541	2546	2554	rBV2	220774	378151	1.06%	0.133%
52	18.027	2554	2557	2560	rVB	39830	48104	0.14%	0.017%
53	18.262	2593	2597	2603	rVB4	66693	110030	0.31%	0.039%
54	18.904	2703	2706	2710	rBV	40310	54824	0.15%	0.019%
55	19.068	2731	2734	2738	rBV	70680	95613	0.27%	0.034%
56	19.251	2761	2765	2769	rBV3	98897	140452	0.39%	0.049%
57	19.345	2774	2781	2786	rVV2	108779	223629	0.63%	0.078%
58	19.439	2791	2797	2799	rVV	5580689	8213107	23.08%	2.879%
59	19.468	2799	2802	2810	rVB	11650181	14293620	40.17%	5.010%
60	20.392	2954	2959	2967	rBV	8417503	9672856	27.19%	3.391%
61	20.980	3055	3059	3064	rBV3	644899	914654	2.57%	0.321%
62	21.180	3088	3093	3097	rBV	10218797	10871195	30.56%	3.811%
63	21.233	3097	3102	3111	rVB	4325560	5474031	15.39%	1.919%
64	21.427	3131	3135	3140	rBV	431210	516027	1.45%	0.181%
65	21.856	3204	3208	3214	rBV	901187	1158961	3.26%	0.406%
66	22.315	3282	3286	3292	rVB	1389964	1776373	4.99%	0.623%
67	22.821	3368	3372	3379	rVB	1688224	2460962	6.92%	0.863%
68	23.297	3447	3453	3461	rVB	4779414	8579696	24.11%	3.007%
69	23.386	3464	3468	3472	rVV	1698875	2955376	8.31%	1.036%
70	23.439	3472	3477	3488	rVB	3269143	6027401	16.94%	2.113%
71	24.033	3573	3578	3586	rVB	1566291	2887300	8.12%	1.012%

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72	24.774	3699	3704	3714	rVB	1144322	2531864	7.12%	0.887%
73	25.668	3844	3856	3868	rVB2	12020941	35578650	100.00%	12.471%
74	26.327	3959	3968	3983	rVB	4739753	14195621	39.90%	4.976%

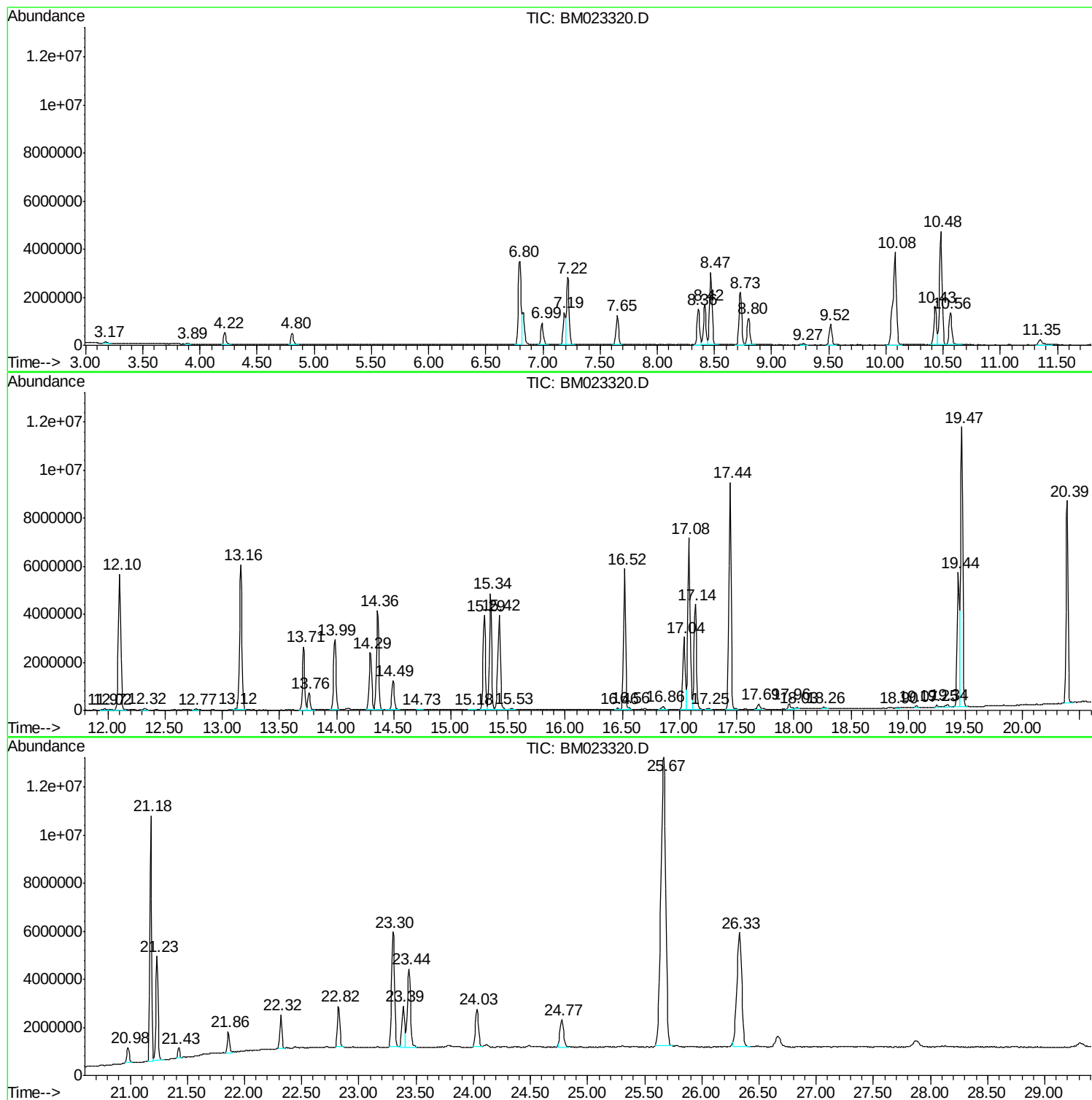
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION

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



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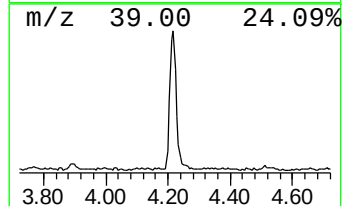
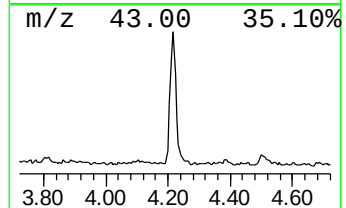
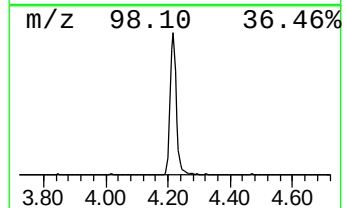
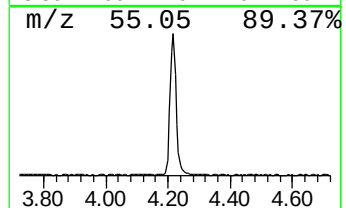
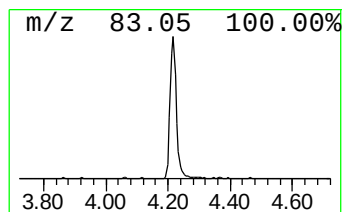
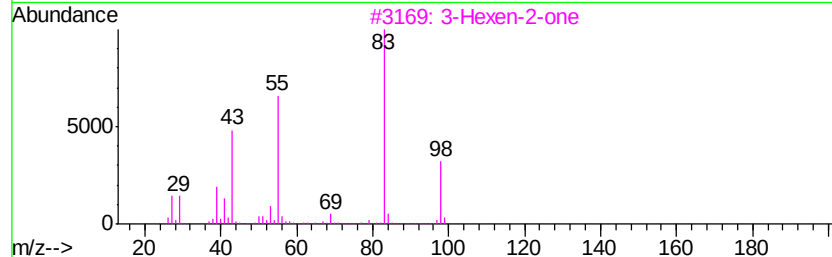
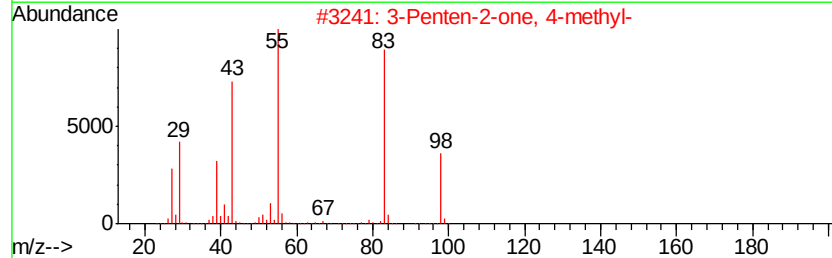
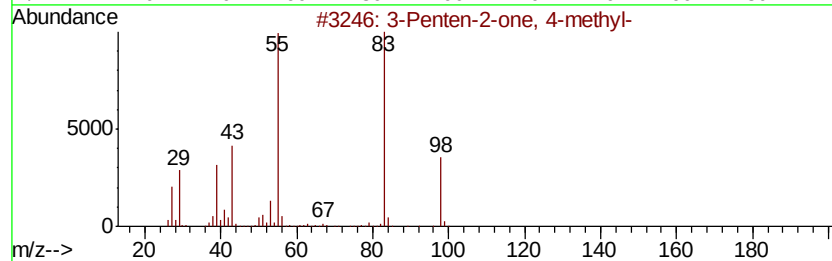
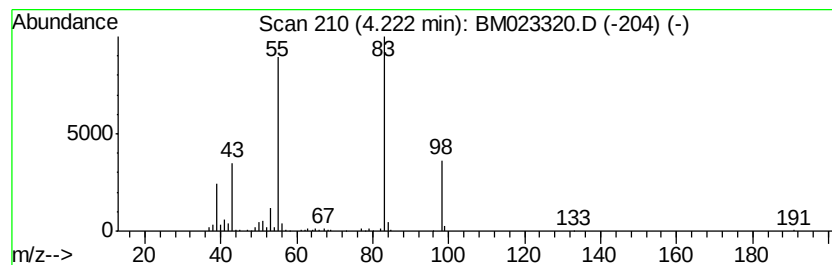
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.22	6.67 ng/ul	650069	1,4-Dichlorobenzene-d4	7.65

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			3-Hexen-2-one	98	C6H10O	000763-93-9	91
5			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90



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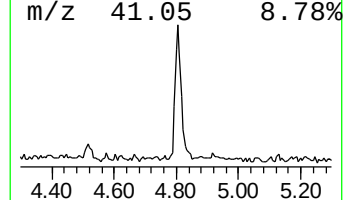
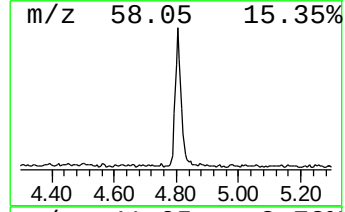
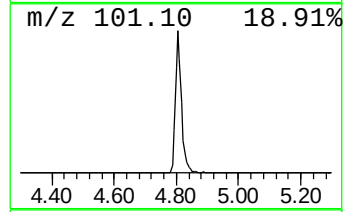
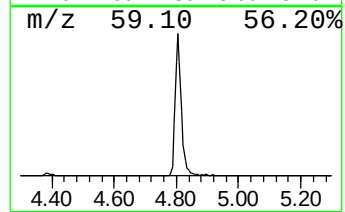
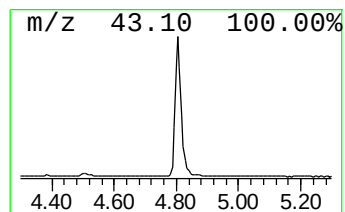
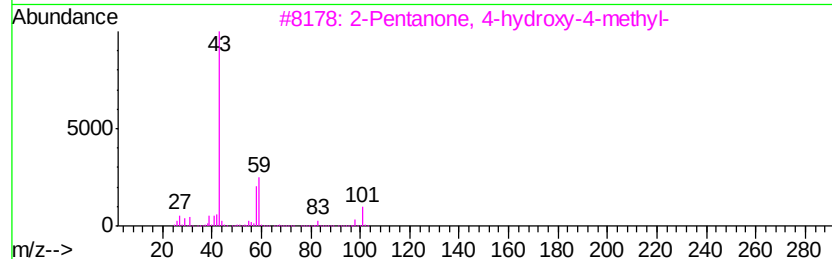
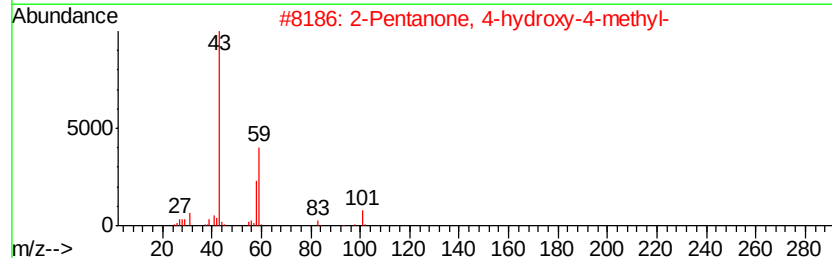
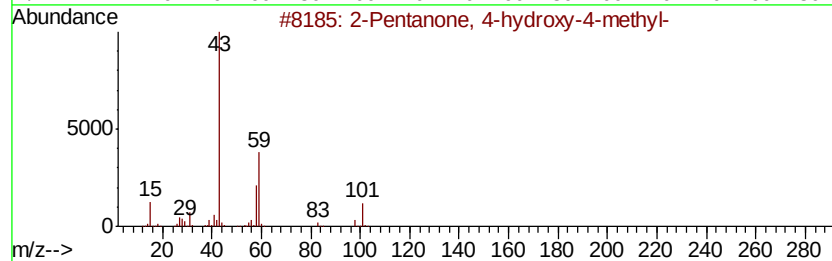
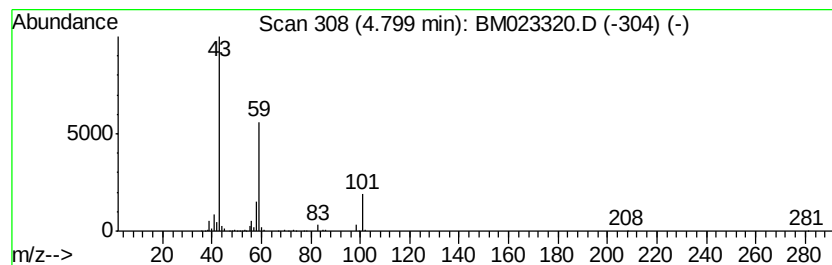
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	6.81 ng/ul	663767	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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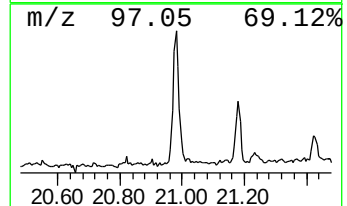
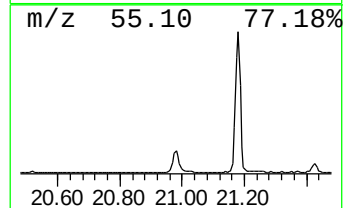
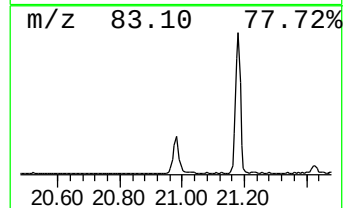
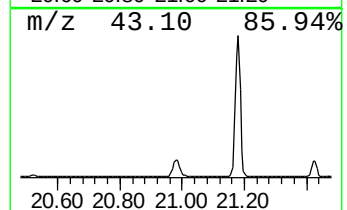
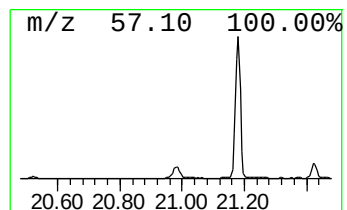
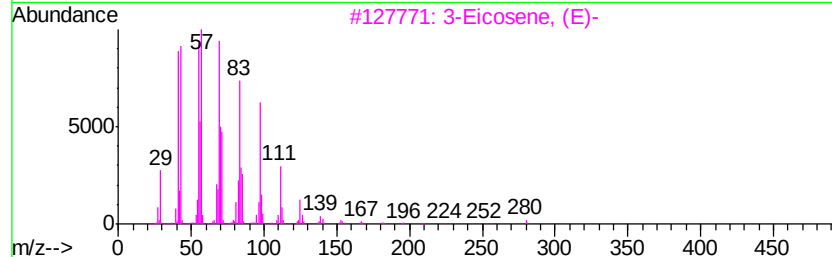
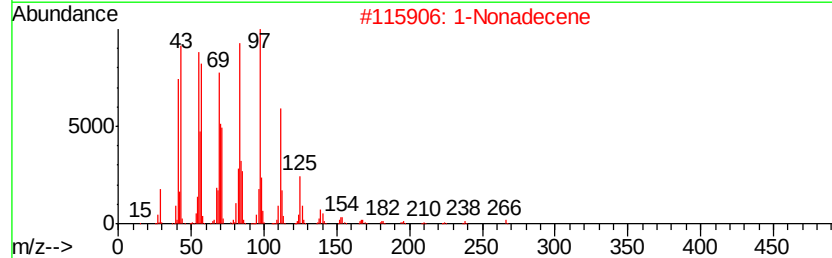
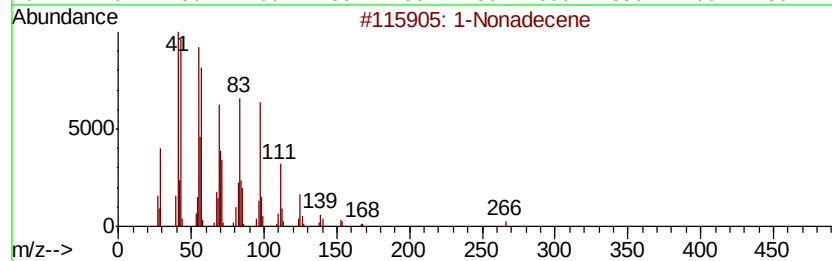
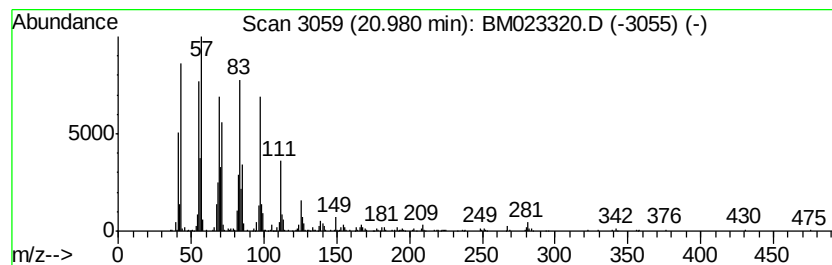
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 (DEL) Alkane: Cyclic20.98 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.98	3.34 ng/ul	914654	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	98
2		1-Nonadecene	266	C19H38	018435-45-5	97
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	96
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	93
5		Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91



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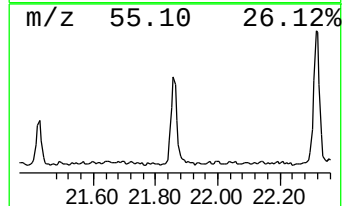
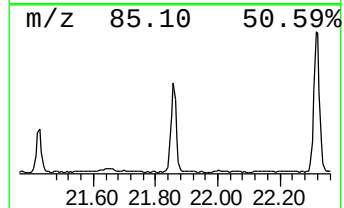
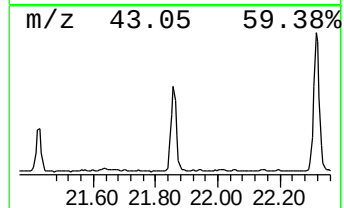
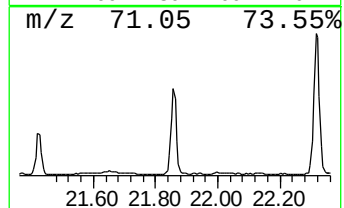
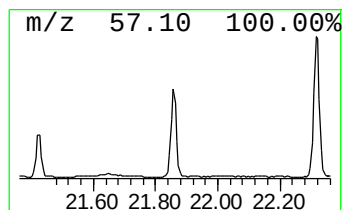
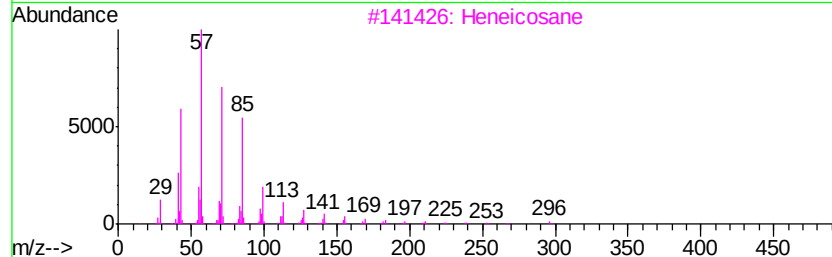
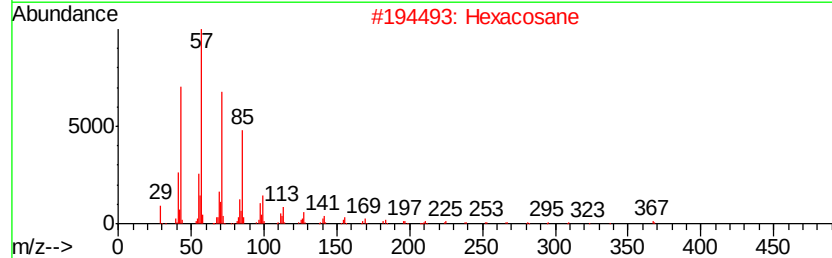
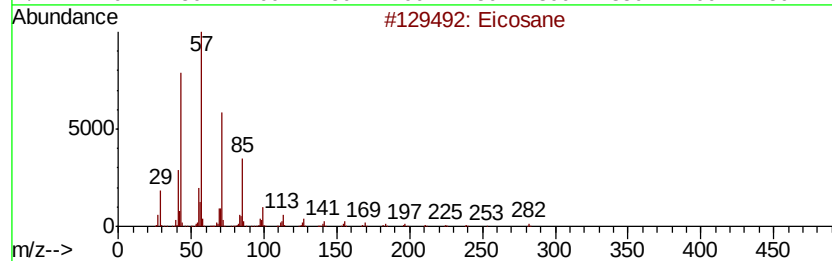
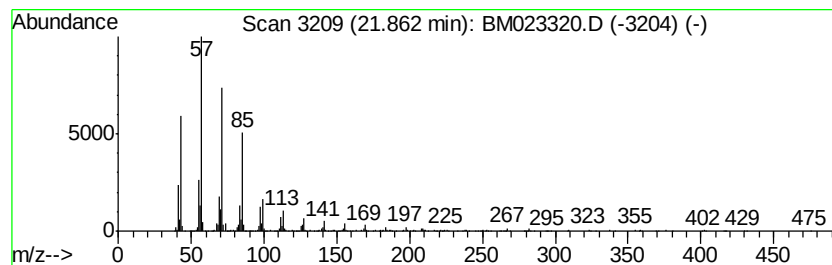
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.86	4.23 ng/ul	1158960	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
X2J25

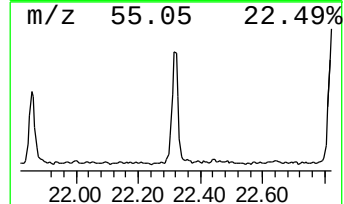
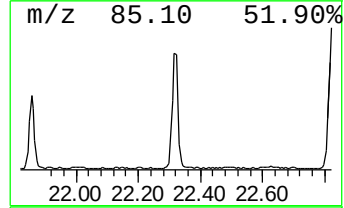
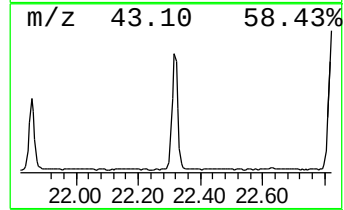
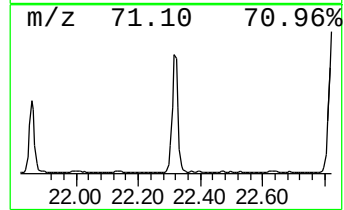
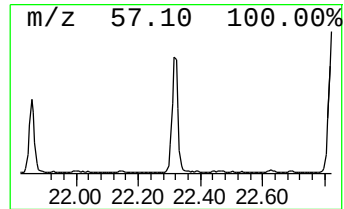
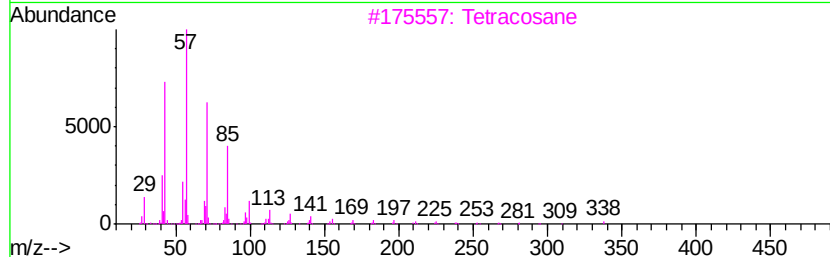
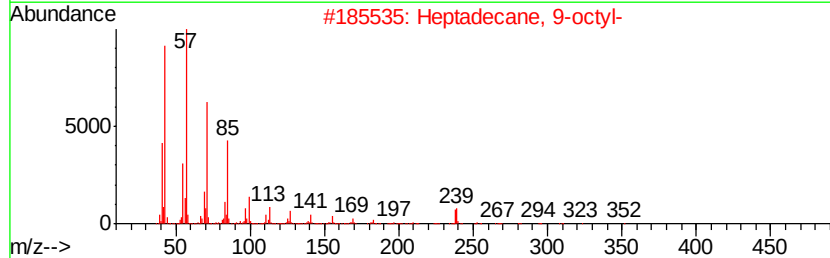
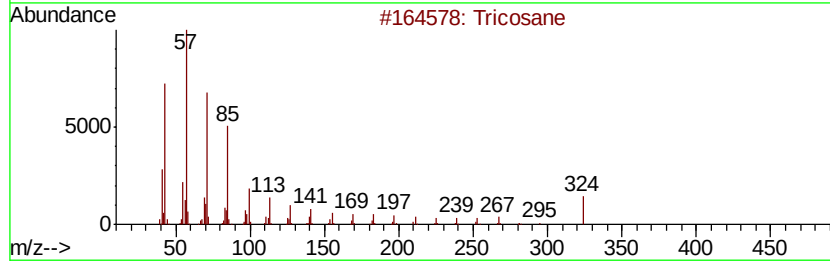
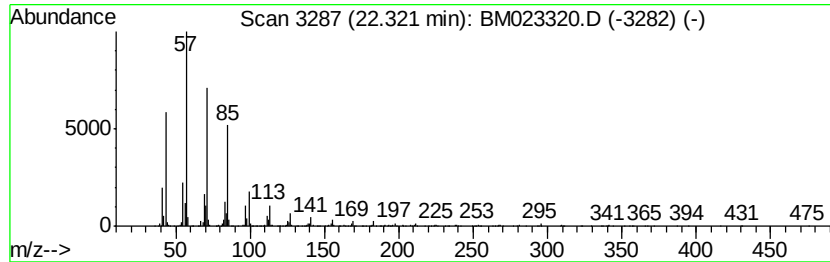
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.32	6.49 ng/ul	1776370	Chrysene-d12	21.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	95
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Octacosane	394	C28H58	000630-02-4	91
5	Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
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Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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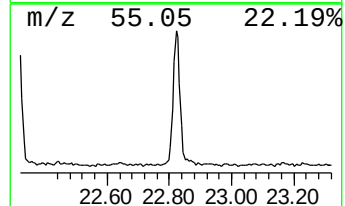
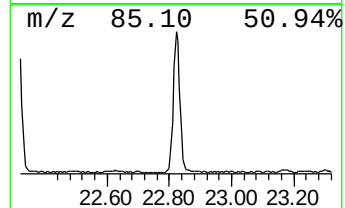
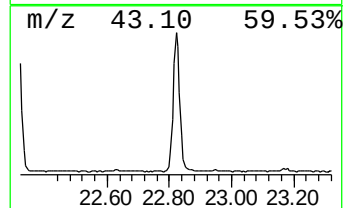
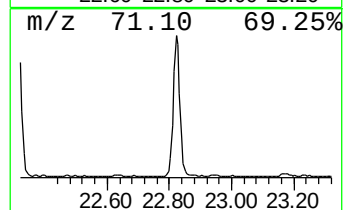
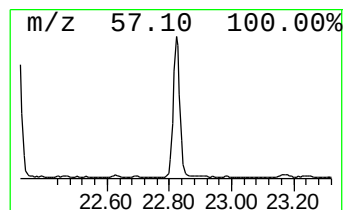
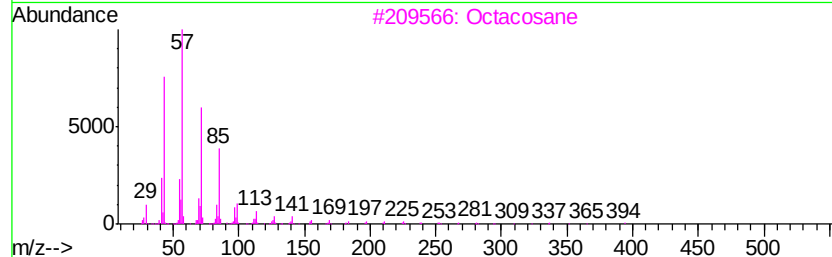
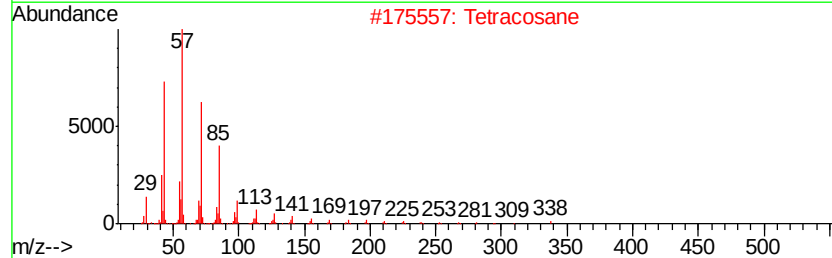
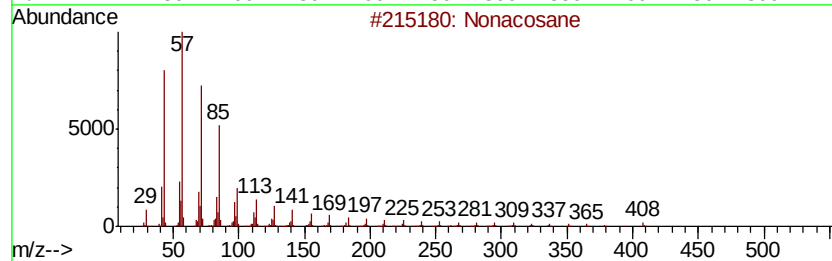
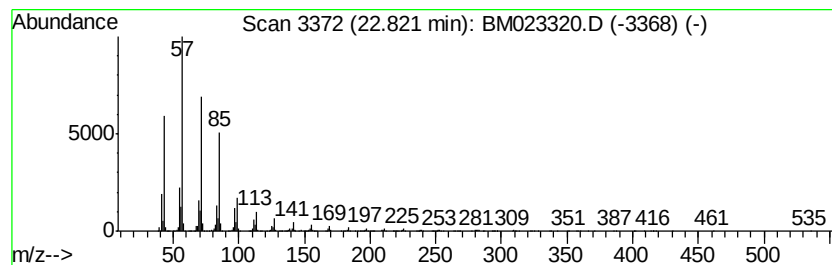
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.82	8.17 ng/ul	2460960	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonacosane	408	C29H60	000630-03-5	93
2		Tetracosane	338	C24H50	000646-31-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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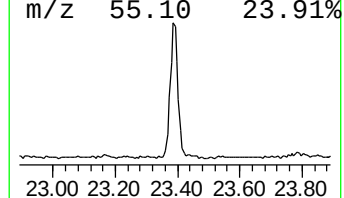
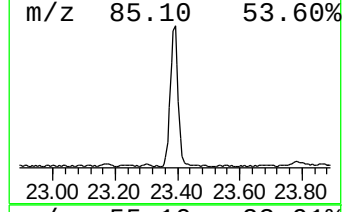
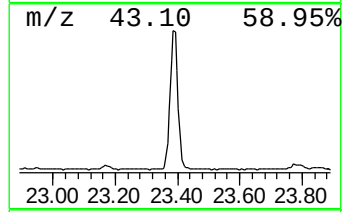
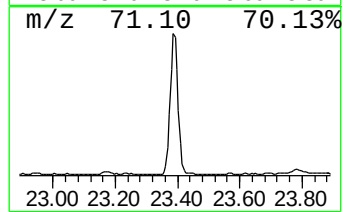
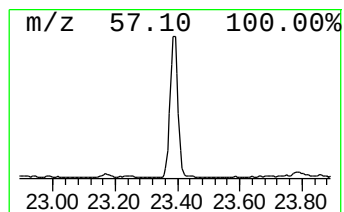
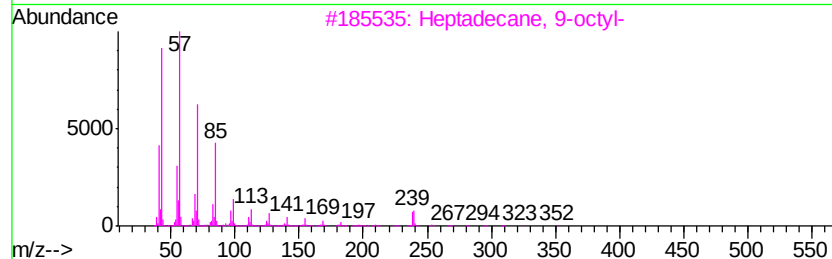
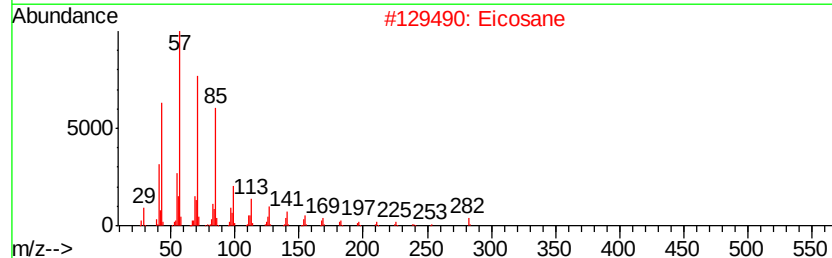
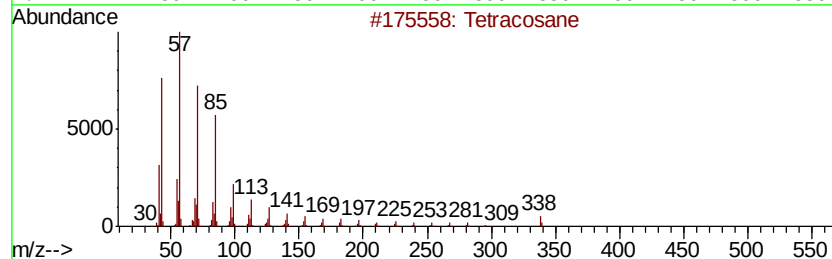
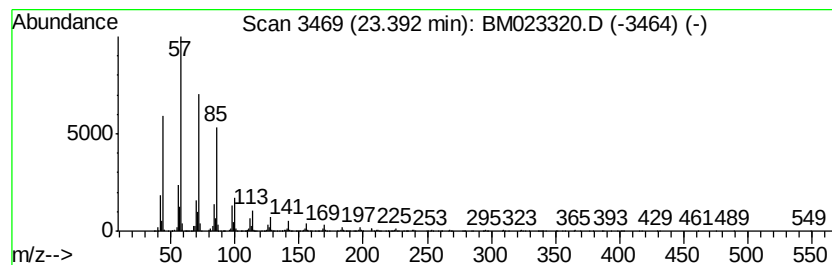
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.39	9.81 ng/ul	2955380	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	98
2		Eicosane	282	C20H42	000112-95-8	97
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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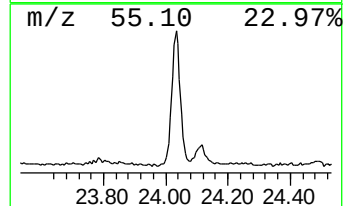
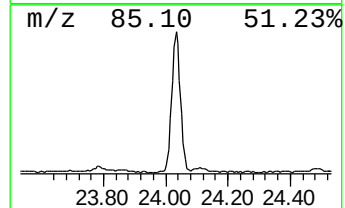
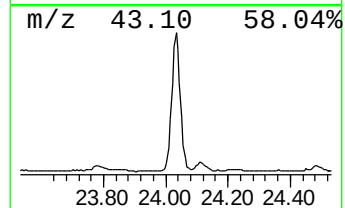
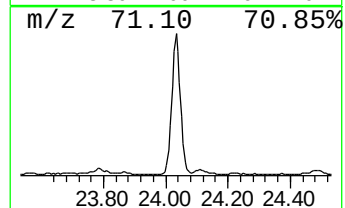
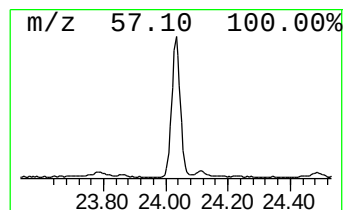
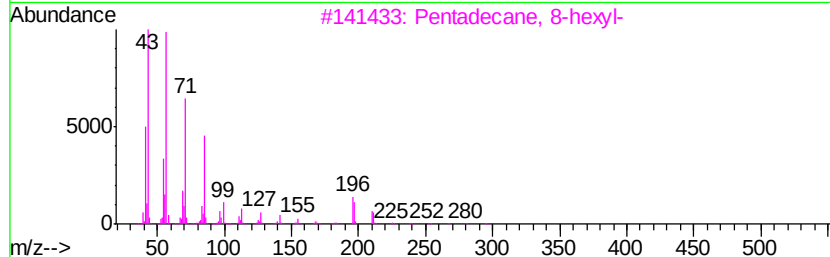
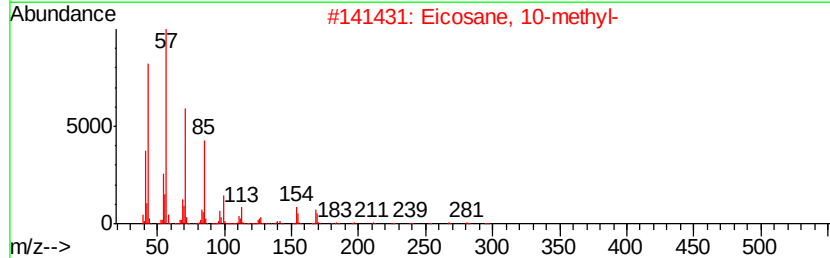
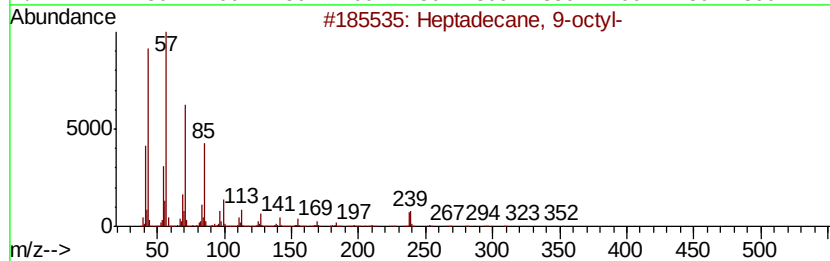
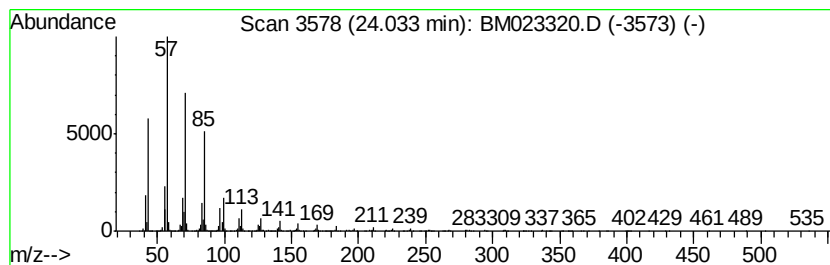
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.03	9.58 ng/ul	2887300	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
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Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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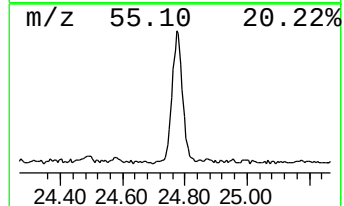
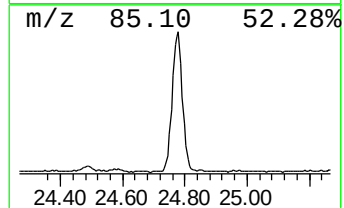
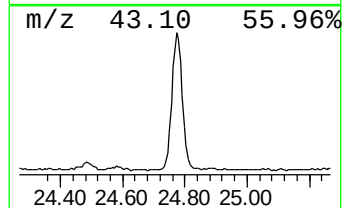
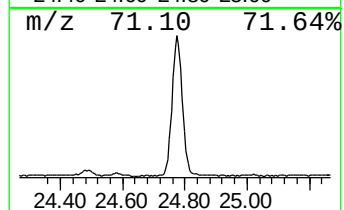
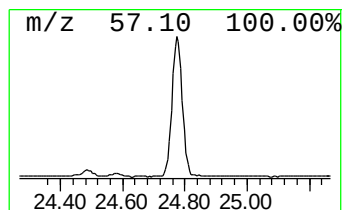
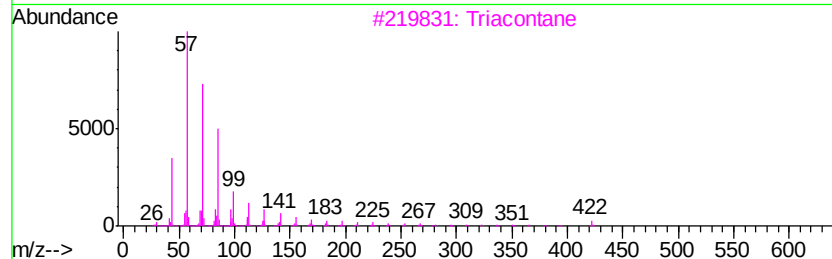
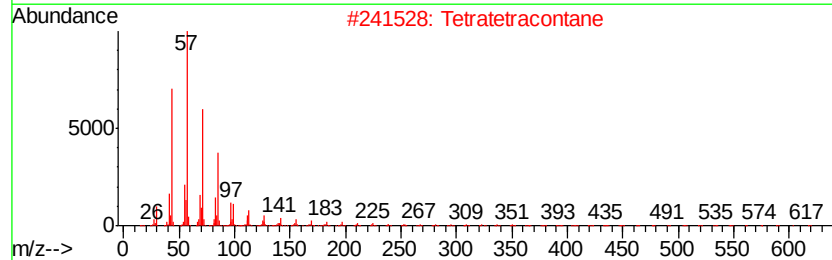
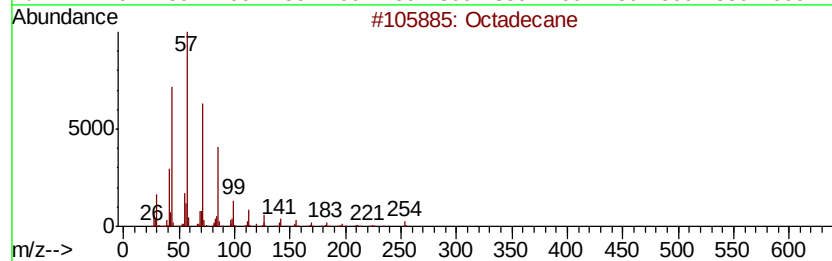
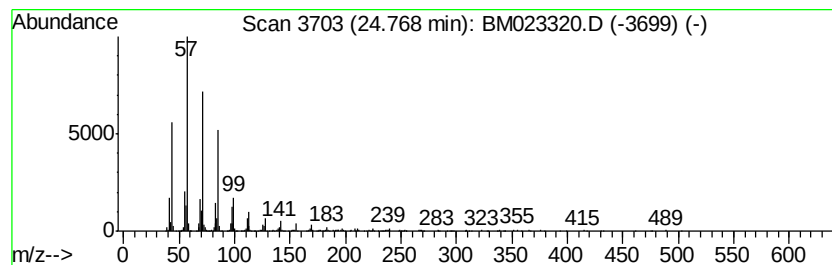
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.77	8.40 ng/ul	2531860	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Hexacosane, 9-octyl-	479	C34H70	055429-83-9	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM102219\
Data File : BM023320.D
Acq On : 22 Oct 2019 12:29
Operator : JU
Sample : K5308-07
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.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
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2-Pentanone, 4-hy...	4.80	6.8	ng/ul	663767	1	7.65	1949470	20.0
(DEL) Alkane: Cyc...	20.98	3.3	ng/ul	914654	5	21.23	5474030	20.0
(DEL) Alkane: Str...	21.86	4.2	ng/ul	1158960	5	21.23	5474030	20.0
(DEL) Alkane: Str...	22.32	6.5	ng/ul	1776370	5	21.23	5474030	20.0
(DEL) Alkane: Str...	22.82	8.2	ng/ul	2460960	6	23.44	6027400	20.0
(DEL) Alkane: Str...	23.39	9.8	ng/ul	2955380	6	23.44	6027400	20.0
(DEL) Alkane: Str...	24.03	9.6	ng/ul	2887300	6	23.44	6027400	20.0
(DEL) Alkane: Str...	24.77	8.4	ng/ul	2531860	6	23.44	6027400	20.0