

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
 Data File : BN009510.D
 Acq On : 31 Jan 2020 14:37
 Operator : CG/JU
 Sample : L1271-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GASZ9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.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	3.081	13	16	24	rVB	78155	109842	3.13%	0.233%
2	3.417	70	73	83	rBV7	3807	11186	0.32%	0.024%
3	3.687	115	119	125	rBV6	18888	34083	0.97%	0.072%
4	3.775	130	134	137	rBV5	7716	9499	0.27%	0.020%
5	3.981	168	169	176	rVB5	3609	5260	0.15%	0.011%
6	4.134	190	195	196	rBV5	3385	4423	0.13%	0.009%
7	4.652	280	283	290	rBV5	6134	10743	0.31%	0.023%
8	5.652	448	453	454	rBV4	3171	5089	0.14%	0.011%
9	5.834	482	484	489	rVB4	2649	4316	0.12%	0.009%
10	6.169	535	541	542	rBV5	6795	9419	0.27%	0.020%
11	6.599	610	614	617	rBV3	2494	3642	0.10%	0.008%
12	6.652	617	623	639	rVV	745533	1304047	37.11%	2.770%
13	6.805	644	649	664	rVB	543049	873331	24.85%	1.855%
14	6.993	675	681	695	rVB	770444	1179634	33.57%	2.506%
15	7.452	753	759	769	rVV	639565	977543	27.82%	2.077%
16	7.540	769	774	780	rVB5	11391	17097	0.49%	0.036%
17	8.163	873	880	892	rBV	917205	1431961	40.75%	3.042%
18	8.593	946	953	964	rBV	741050	1199285	34.13%	2.548%
19	8.740	973	978	981	rBV6	3417	5687	0.16%	0.012%
20	9.052	1027	1031	1035	rVB3	7010	9172	0.26%	0.019%
21	9.305	1067	1074	1087	rBV	634160	1016023	28.91%	2.158%
22	9.840	1158	1165	1177	rBV	845796	1452145	41.32%	3.085%
23	10.116	1207	1212	1214	rBV3	3141	4421	0.13%	0.009%
24	10.204	1218	1227	1234	rBV	870182	1427597	40.63%	3.033%
25	10.352	1245	1252	1269	rBV	747369	1349477	38.40%	2.867%
26	11.157	1386	1389	1392	rBV	2295	3565	0.10%	0.008%
27	11.804	1493	1499	1505	rBV2	13410	21554	0.61%	0.046%
28	11.881	1510	1512	1516	rVB3	3486	4202	0.12%	0.009%
29	12.934	1684	1691	1693	rBV3	3511	5130	0.15%	0.011%
30	13.004	1698	1703	1708	rBV	17411	25507	0.73%	0.054%
31	13.510	1783	1789	1794	rBV	1532519	2044543	58.18%	4.343%
32	13.557	1794	1797	1807	rVB	351692	493760	14.05%	1.049%
33	13.775	1826	1834	1842	rBV	1768348	2495413	71.01%	5.301%
34	14.028	1873	1877	1881	rVV3	5381	5961	0.17%	0.013%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.M
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35	14.087	1881	1887	1899	rVB	1302248	1882147	53.56%	3.998%
36	14.316	1920	1926	1945	rBV	545339	1019834	29.02%	2.166%
37	14.469	1949	1952	1959	rVV7	4352	8930	0.25%	0.019%
38	14.540	1959	1964	1970	rVB8	8714	16175	0.46%	0.034%
39	14.645	1979	1982	1988	rBV5	3388	5351	0.15%	0.011%
40	14.928	2025	2030	2035	rBV5	4753	9197	0.26%	0.020%
41	14.981	2035	2039	2046	rVB4	8433	12394	0.35%	0.026%
42	15.092	2051	2058	2064	rBV	2141425	3021457	85.98%	6.419%
43	15.228	2076	2081	2094	rBV	648999	1000165	28.46%	2.125%
44	15.334	2094	2099	2104	rVB3	21730	30627	0.87%	0.065%
45	15.545	2131	2135	2137	rVB5	4728	5455	0.16%	0.012%
46	15.787	2171	2176	2177	rBV4	2677	4209	0.12%	0.009%
47	15.822	2178	2182	2184	rVV3	7358	10750	0.31%	0.023%
48	15.839	2184	2185	2189	rVV3	6615	9485	0.27%	0.020%
49	15.881	2189	2192	2195	rVV3	8034	10948	0.31%	0.023%
50	15.945	2198	2203	2204	rBV5	3586	4325	0.12%	0.009%
51	16.104	2225	2230	2233	rBV4	3687	5497	0.16%	0.012%
52	16.286	2257	2261	2265	rVB3	6958	9248	0.26%	0.020%
53	16.592	2309	2313	2316	rBV5	3453	4808	0.14%	0.010%
54	16.639	2316	2321	2325	rVB4	10672	15244	0.43%	0.032%
55	16.839	2349	2355	2362	rVV	1591713	2149879	61.18%	4.567%
56	16.939	2366	2372	2385	rVV	2302825	3150321	89.65%	6.692%
57	17.057	2388	2392	2393	rBV3	4347	5402	0.15%	0.011%
58	17.181	2410	2413	2417	rVB5	5607	8108	0.23%	0.017%
59	17.257	2419	2426	2430	rBV3	44688	63650	1.81%	0.135%
60	17.645	2488	2492	2494	rBV4	4482	6036	0.17%	0.013%
61	17.733	2502	2507	2509	rBV3	4163	5818	0.17%	0.012%
62	17.775	2509	2514	2523	rBV	313126	501791	14.28%	1.066%
63	17.957	2543	2545	2548	rVB4	4304	4098	0.12%	0.009%
64	18.204	2582	2587	2591	rBV2	49038	53974	1.54%	0.115%
65	18.345	2609	2611	2615	rVB4	3987	4380	0.12%	0.009%
66	18.445	2626	2628	2631	rBV2	4650	7212	0.21%	0.015%
67	18.586	2647	2652	2653	rVB4	5985	8550	0.24%	0.018%
68	18.639	2656	2661	2670	rBV	170460	241582	6.87%	0.513%
69	18.775	2682	2684	2687	rVB4	6721	6939	0.20%	0.015%
70	18.880	2698	2702	2705	rBV2	29665	36297	1.03%	0.077%
71	18.939	2710	2712	2713	rVV2	10106	8901	0.25%	0.019%

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72	18.963	2714	2716	2720	rVB5	28565	29616	0.84%	0.063%
73	19.004	2720	2723	2726	rVB4	11672	13211	0.38%	0.028%
74	19.063	2729	2733	2741	rBV	104105	173333	4.93%	0.368%
75	19.128	2741	2744	2752	rVB	25730	43249	1.23%	0.092%
76	19.245	2757	2764	2770	rVV	2649658	3513999	100.00%	7.465%
77	19.298	2770	2773	2776	rVB3	19209	20901	0.59%	0.044%
78	19.798	2854	2858	2862	rBV3	89512	123694	3.52%	0.263%
79	19.839	2862	2865	2868	rVB	45874	54330	1.55%	0.115%
80	20.339	2946	2950	2953	rVB	74934	85439	2.43%	0.181%
81	20.792	3023	3027	3034	rBV	973030	1228894	34.97%	2.611%
82	21.051	3065	3071	3080	rBV	1753505	2299239	65.43%	4.884%
83	21.239	3100	3103	3107	rBV	148382	158139	4.50%	0.336%
84	21.657	3171	3174	3180	rBV	226479	317926	9.05%	0.675%
85	22.092	3245	3248	3253	rVB	253293	314503	8.95%	0.668%
86	22.210	3265	3268	3275	rVB	227658	282787	8.05%	0.601%
87	22.569	3325	3329	3334	rBV	357847	455243	12.96%	0.967%
88	23.033	3402	3408	3414	rBV	1830913	3122905	88.87%	6.634%
89	23.098	3414	3419	3423	rVV	333709	508458	14.47%	1.080%
90	23.163	3424	3430	3437	rVB	1264364	2174706	61.89%	4.620%
91	23.686	3515	3519	3525	rVB	292264	501623	14.27%	1.066%
92	23.763	3529	3532	3542	rVB3	196822	342552	9.75%	0.728%
93	24.368	3632	3635	3643	rVB3	217184	405509	11.54%	0.861%

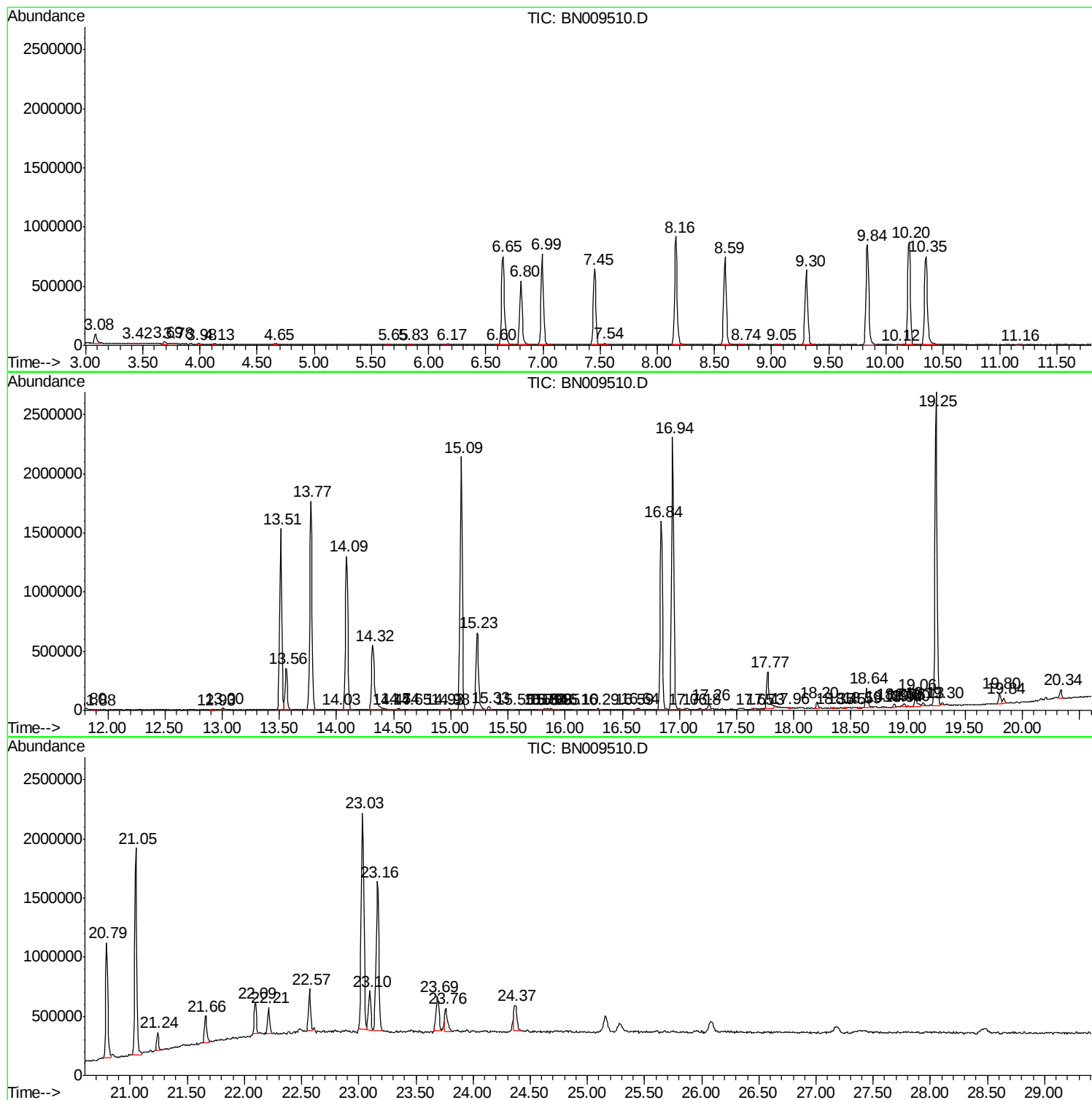
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.M
Quant Title : SVOA CALIBRATION

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



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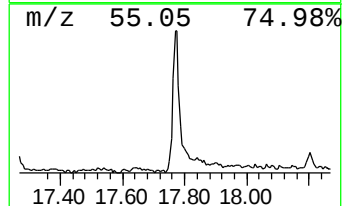
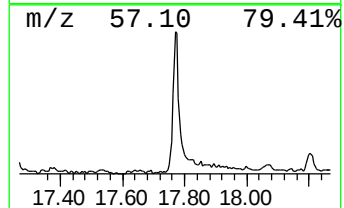
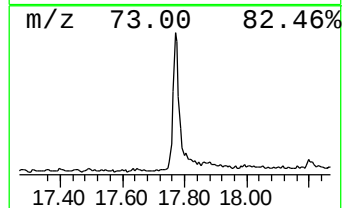
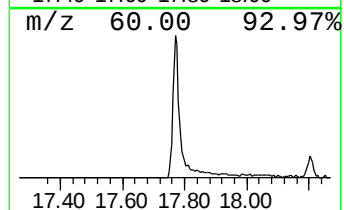
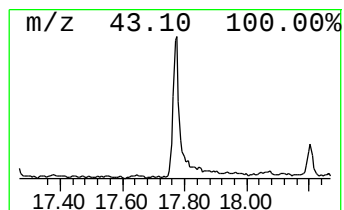
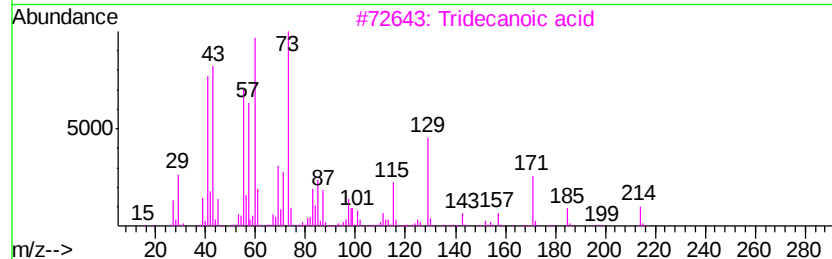
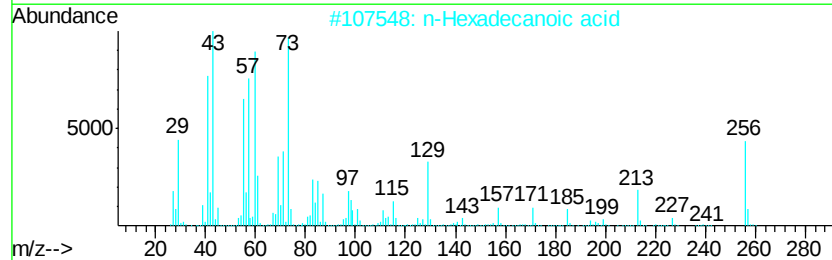
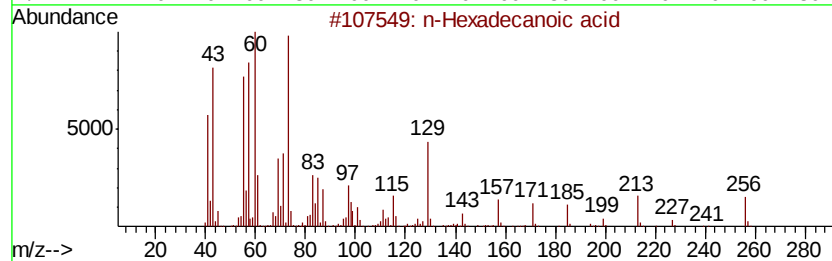
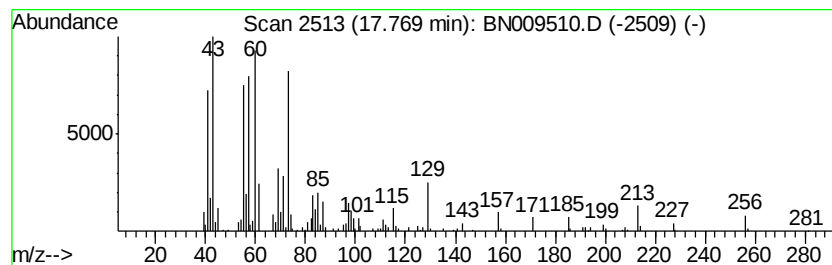
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.77	4.67 ng/ul	501791	Phenanthrene-d10		16.84	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	94
2	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	91
3	Tridecanoic acid		214	C13H26O2	000638-53-9	90
4	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	87
5	Pentadecanoic acid		242	C15H30O2	001002-84-2	87



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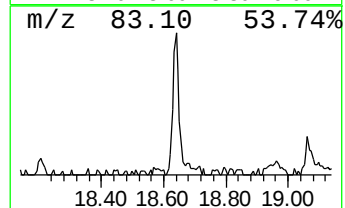
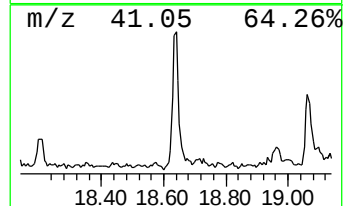
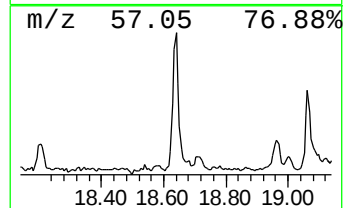
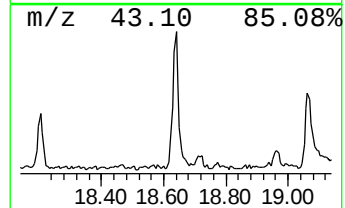
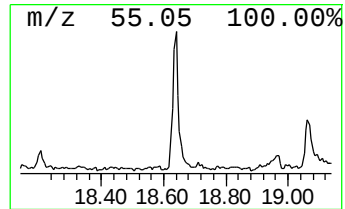
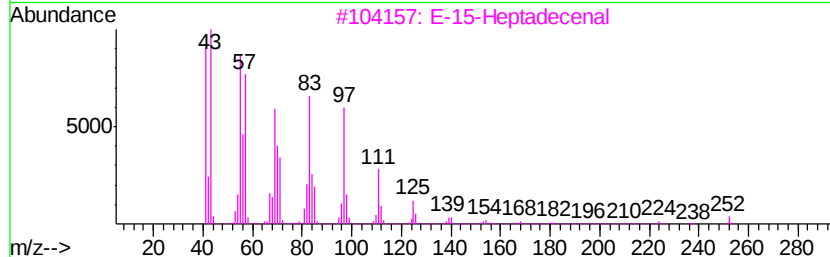
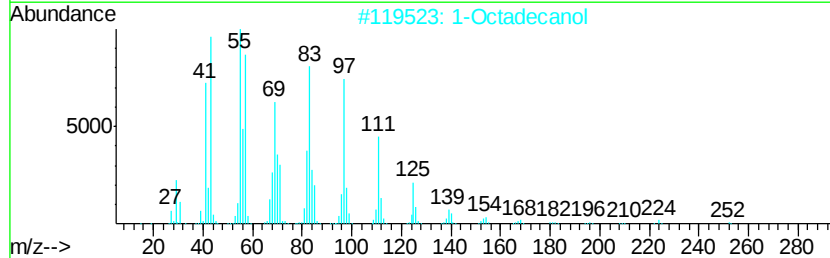
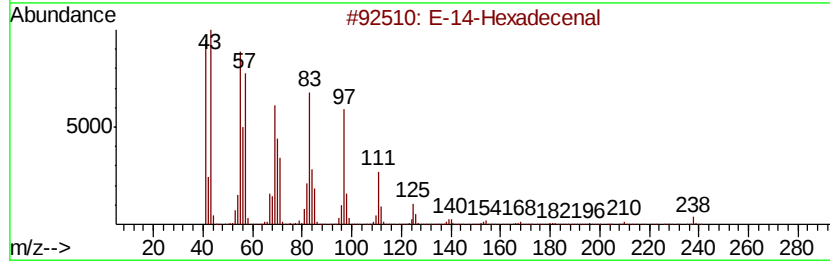
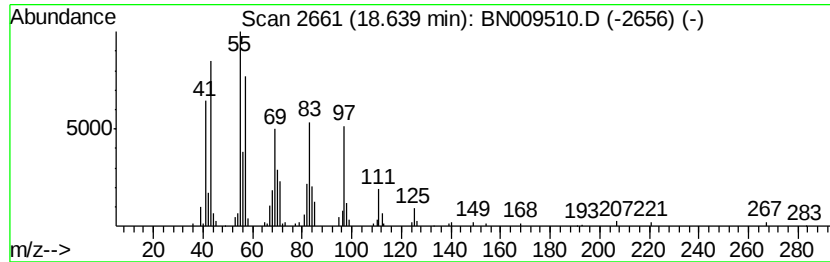
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 E-14-Hexadecenal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
18.64	2.25 ng/ul	241582	Phenanthrene-d10		16.84		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-14-Hexadecenal			238	C16H30O	330207-53-9	91
2	1-Octadecanol			270	C18H38O	000112-92-5	91
3	E-15-Heptadecenal			252	C17H32O	1000130-97-9	91
4	1-Heneicosanol			312	C21H44O	015594-90-8	91
5	n-Nonadecanol-1			284	C19H40O	001454-84-8	90



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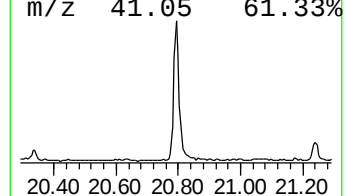
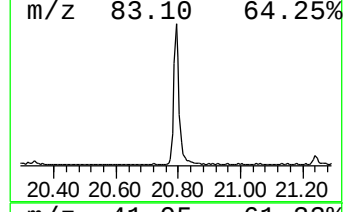
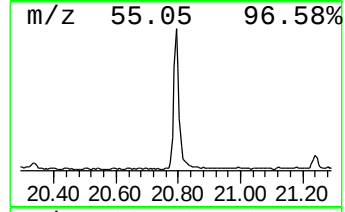
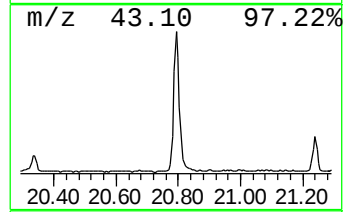
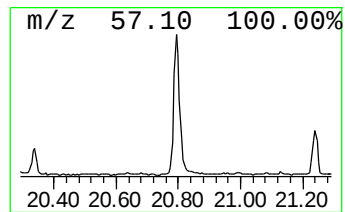
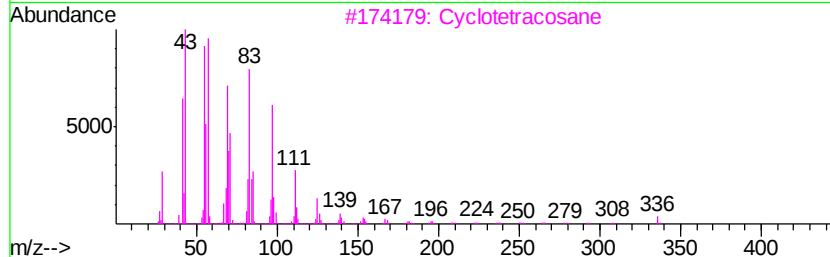
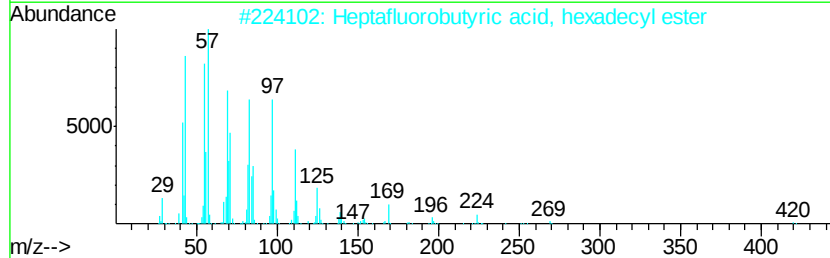
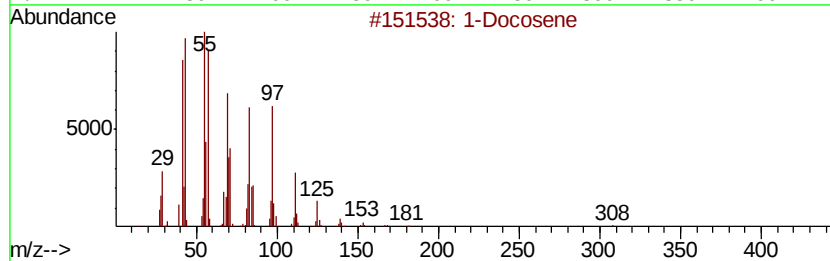
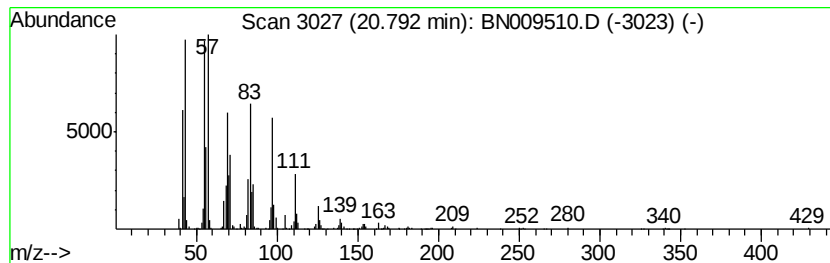
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	10.69 ng/ul	1228890	Chrysene-d12	21.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	95
2		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
3		Cyclotetracosane	336	C24H48	000297-03-0	91
4		Behenic alcohol	326	C22H46O	000661-19-8	91
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
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ClientSampleId :
GASZ9

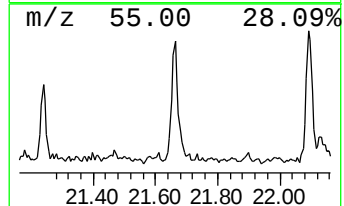
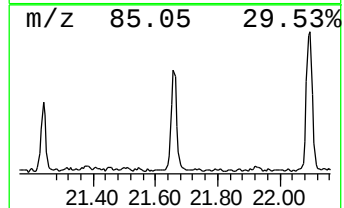
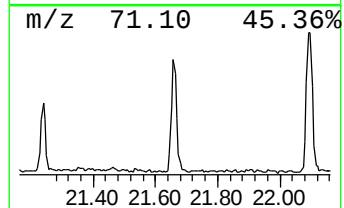
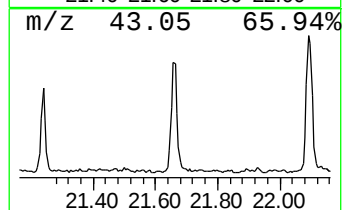
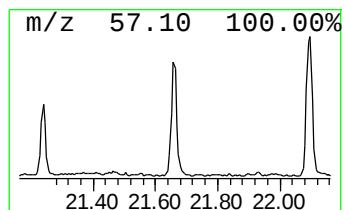
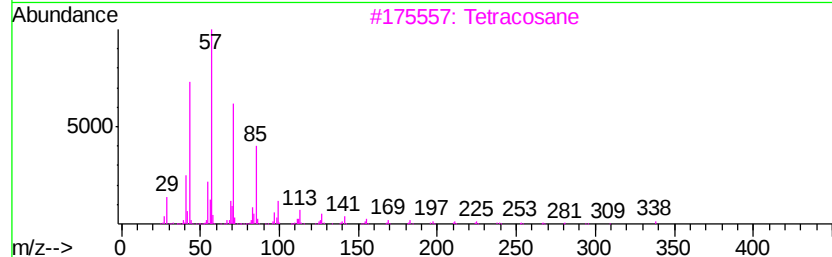
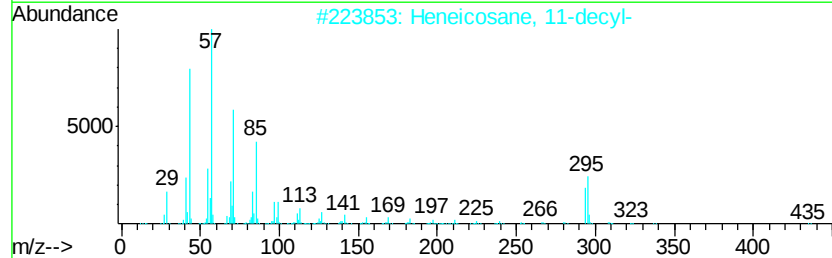
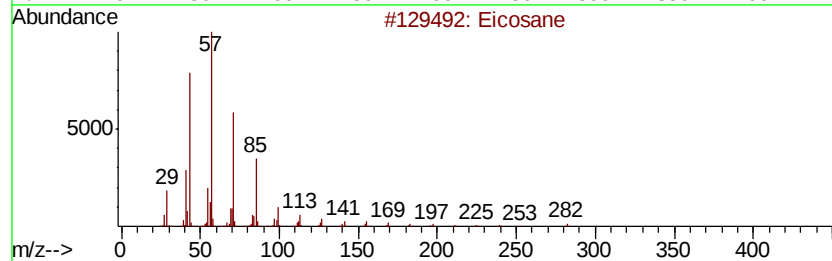
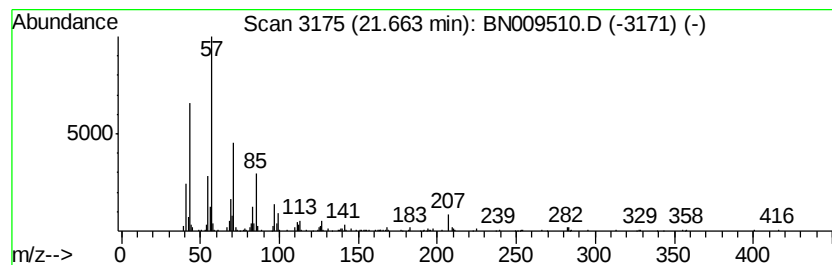
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Quant Title : SVOA CALIBRATION

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.
21.66	2.77 ng/ul	317926	Chrysene-d12	21.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86
3		Tetracosane	338	C24H50	000646-31-1	80
4		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	74
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
Operator : CG/JU
Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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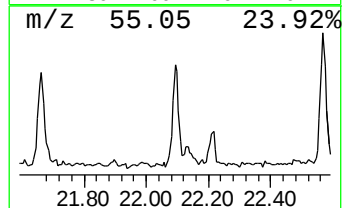
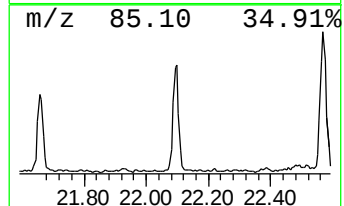
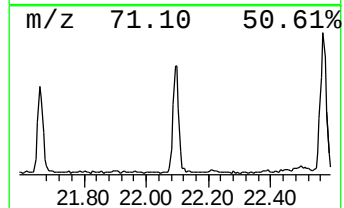
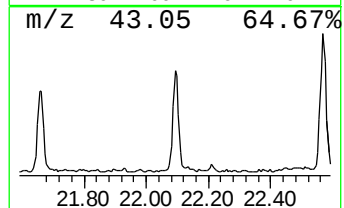
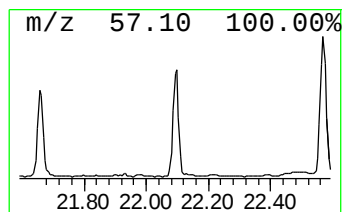
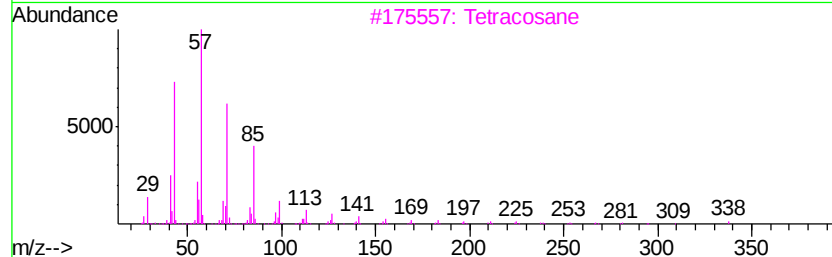
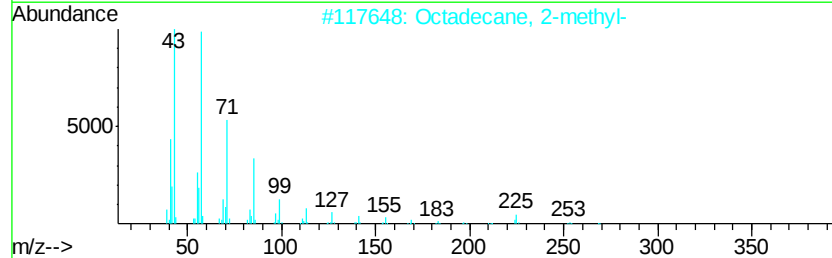
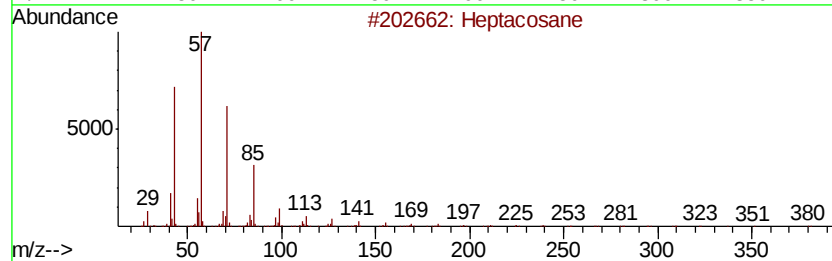
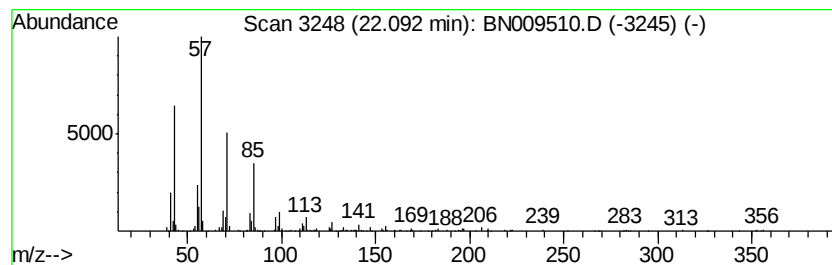
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.09	2.74 ng/ul	314503	Chrysene-d12	21.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	90
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
3		Tetracosane	338	C24H50	000646-31-1	87
4		Hexacosane	366	C26H54	000630-01-3	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
Operator : CG/JU
Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GASZ9

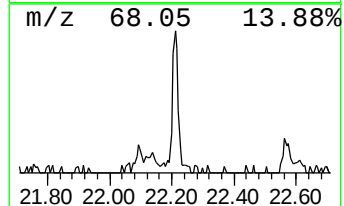
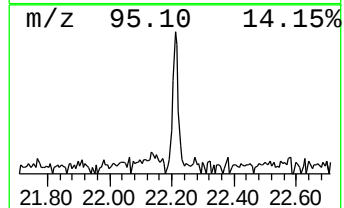
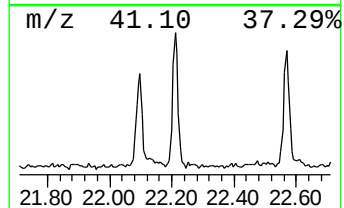
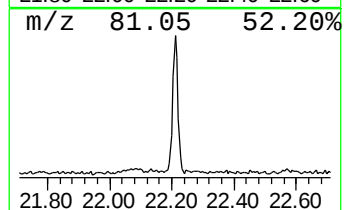
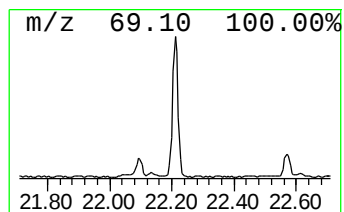
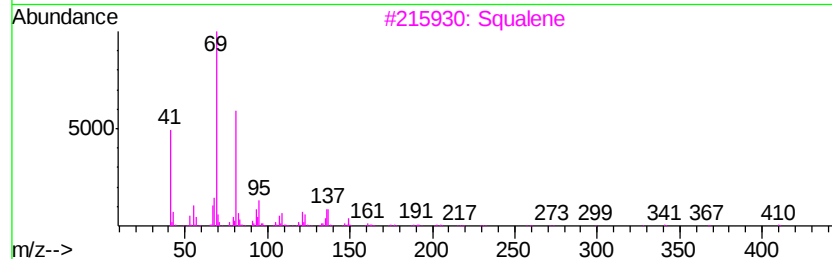
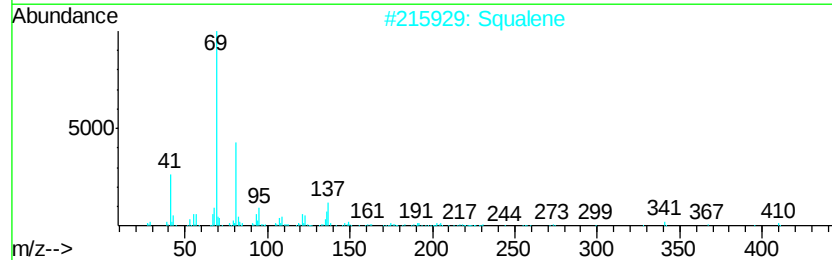
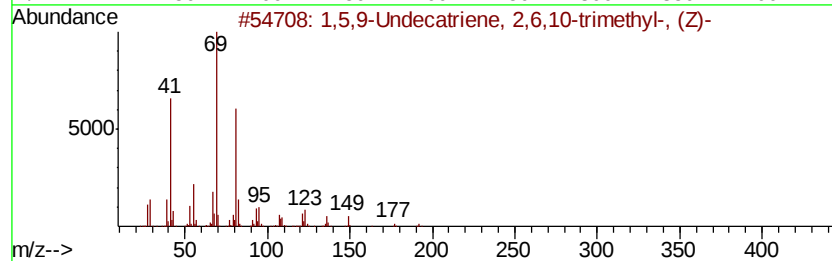
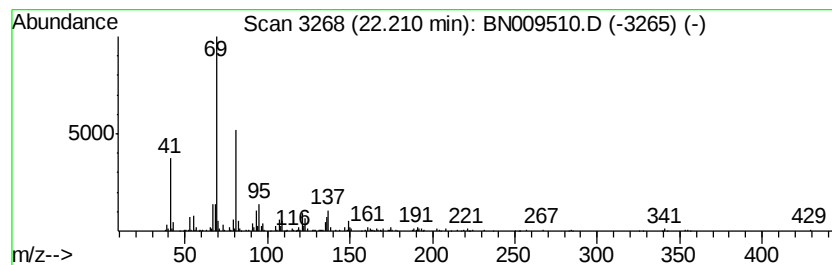
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,5,9-Undecatriene, 2,6,10-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.21	2.60 ng/ul	282787	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	90
2		Squalene	410	C30H50	000111-02-4	87
3		Squalene	410	C30H50	000111-02-4	87
4		Supraene	410	C30H50	007683-64-9	87
5		trans-Farnesol	222	C15H26O	000106-28-5	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
Operator : CG/JU
Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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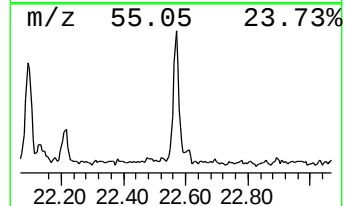
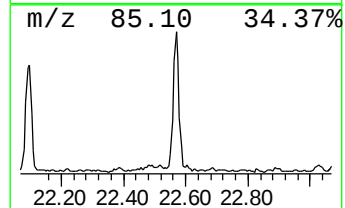
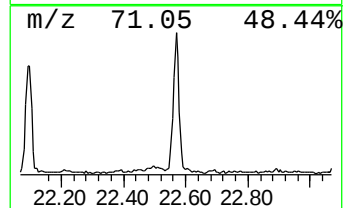
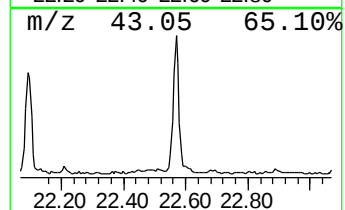
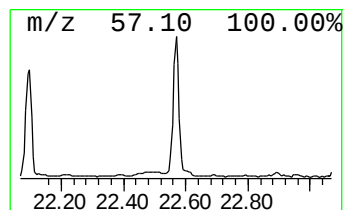
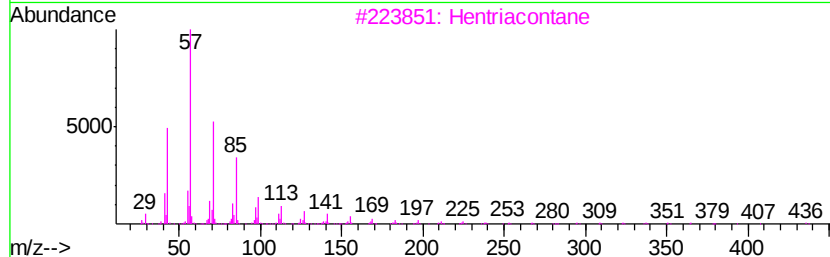
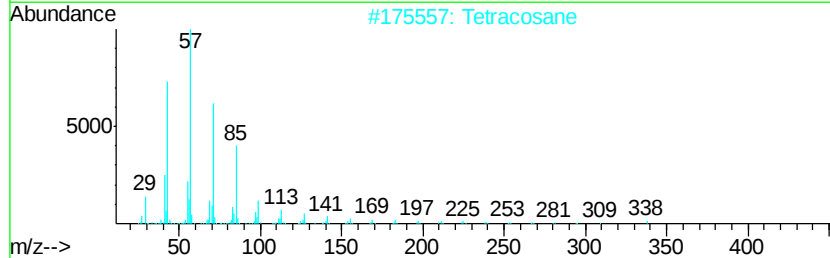
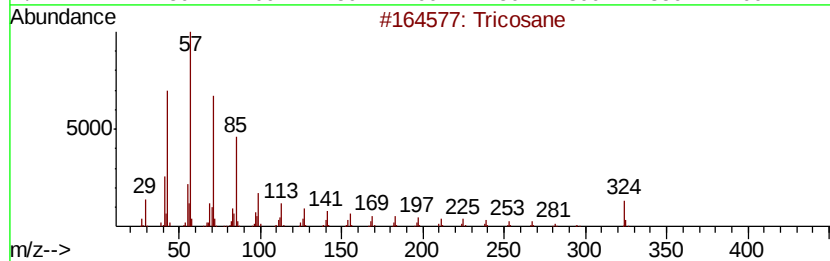
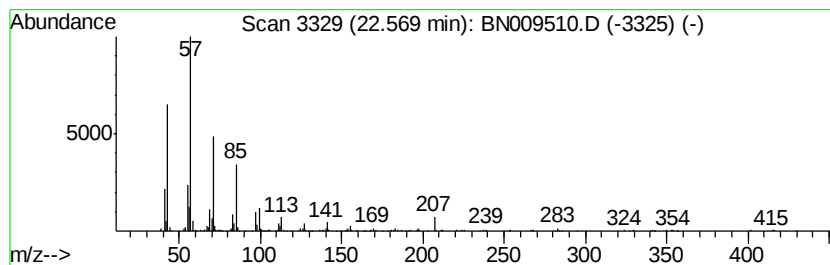
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.57	4.19 ng/ul	455243	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricosane	324	C23H48	000638-67-5	93
2		Tetracosane	338	C24H50	000646-31-1	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
 Data File : BN009510.D
 Acq On : 31 Jan 2020 14:37
 Operator : CG/JU
 Sample : L1271-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
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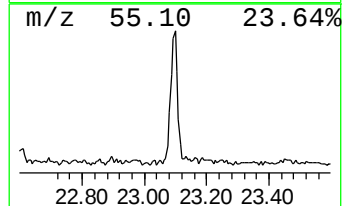
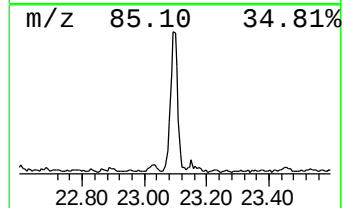
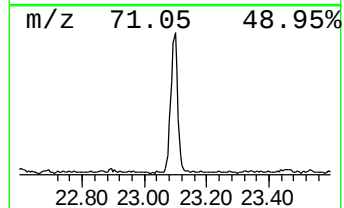
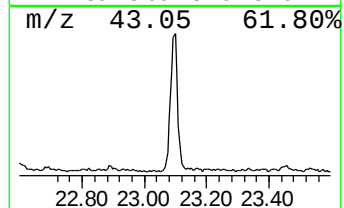
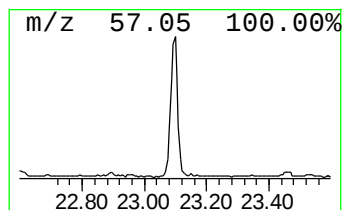
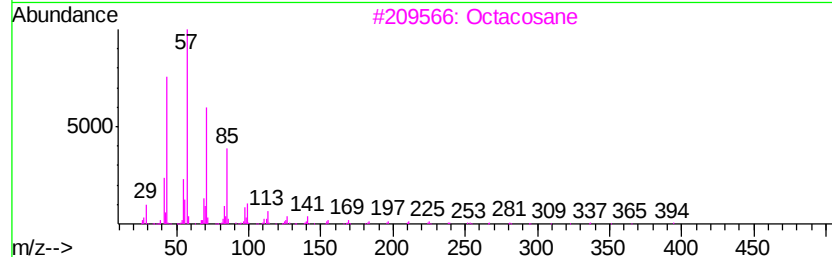
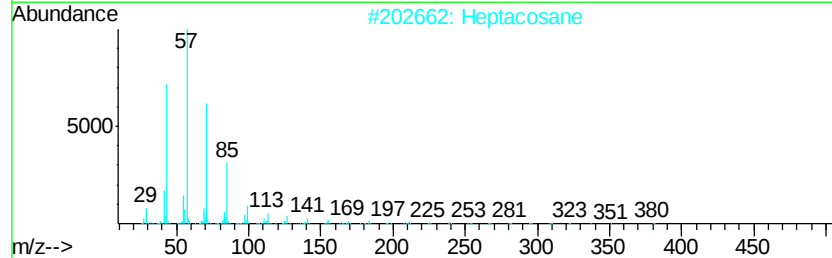
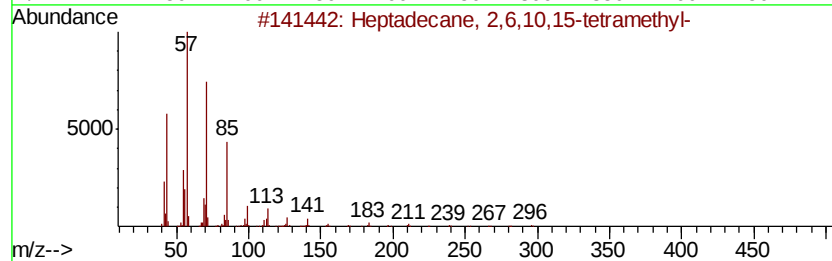
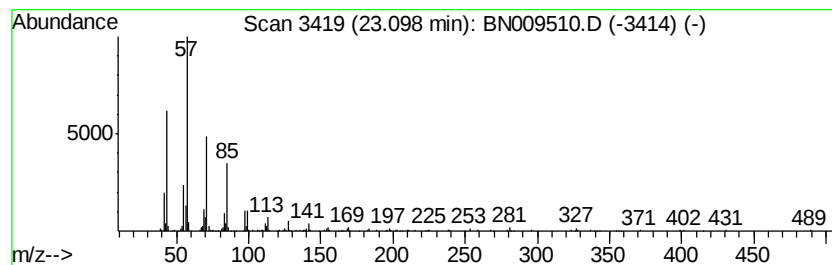
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 Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.10	4.68 ng/ul	508458	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	97
2		Heptacosane	380	C27H56	000593-49-7	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Hentriacontane	437	C31H64	000630-04-6	87
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
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Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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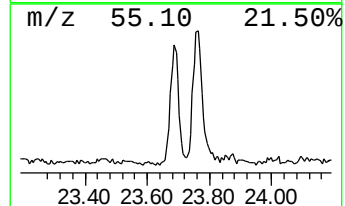
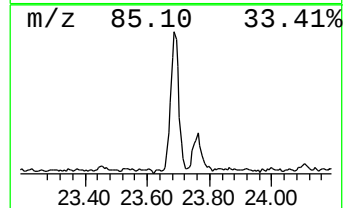
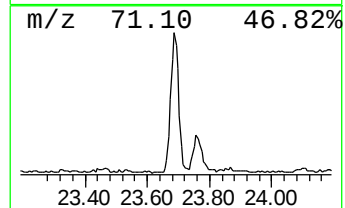
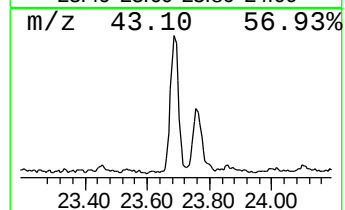
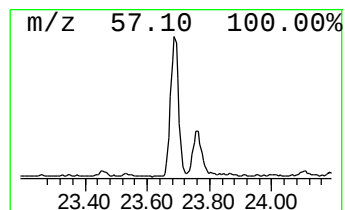
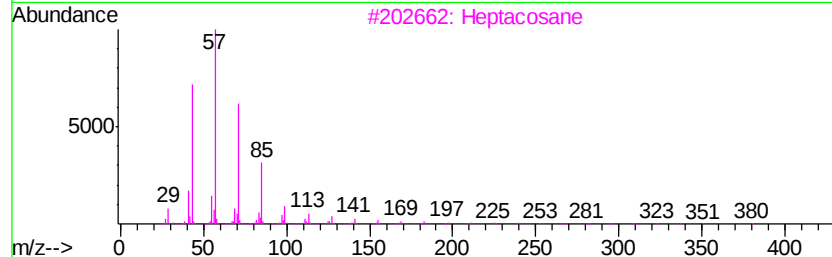
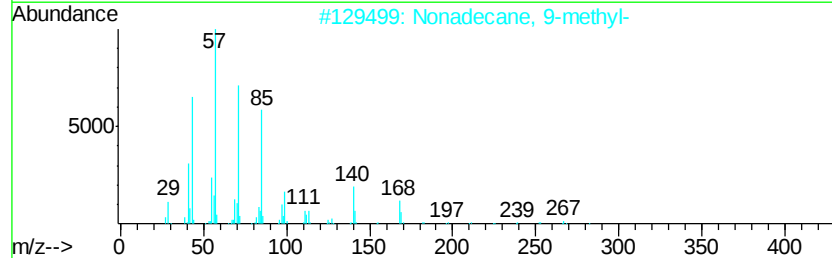
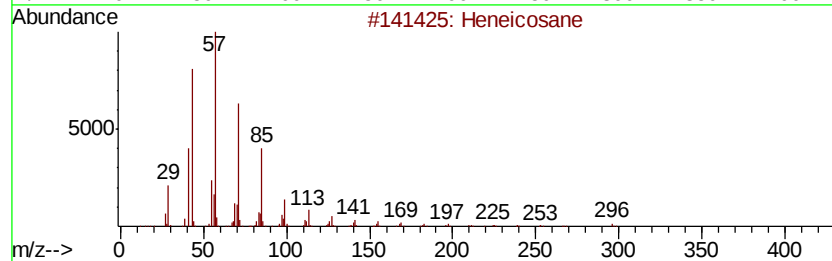
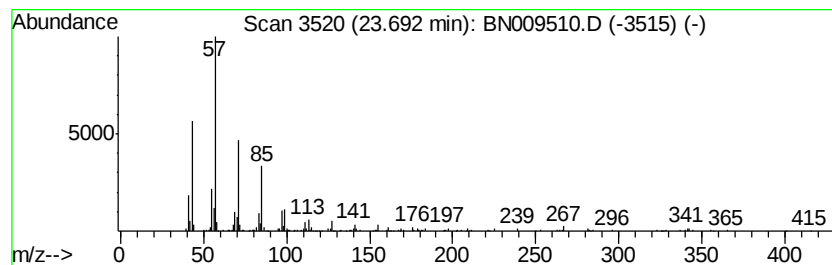
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.69	4.61 ng/ul	501623	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	96
2		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	89
3		Heptacosane	380	C27H56	000593-49-7	87
4		Octacosane	394	C28H58	000630-02-4	87
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
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Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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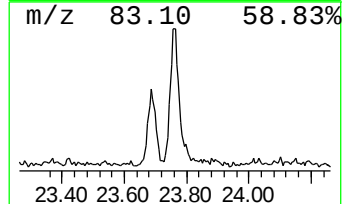
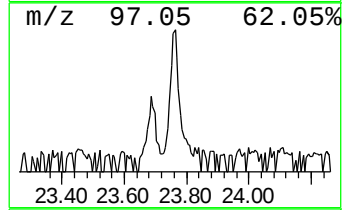
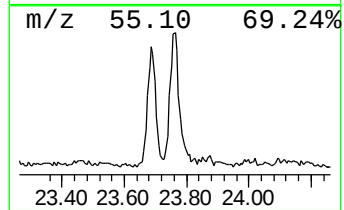
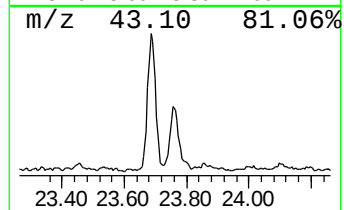
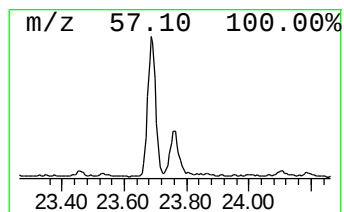
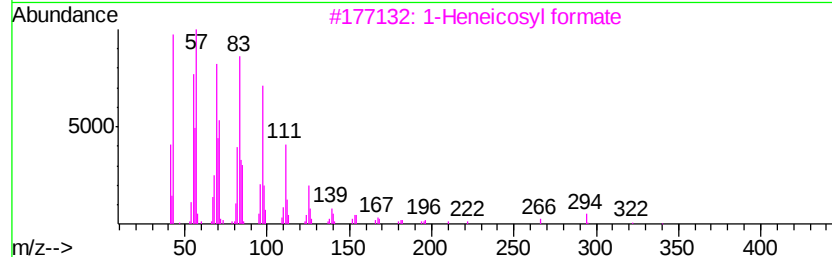
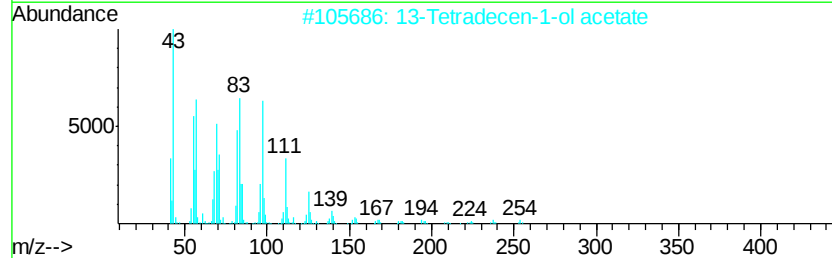
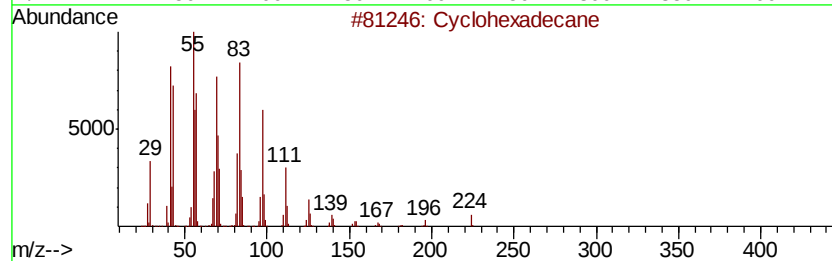
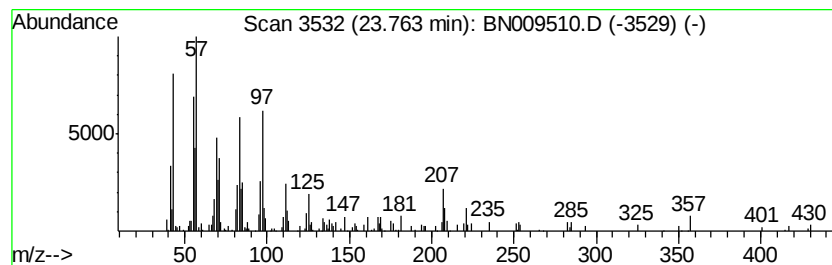
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic23.76 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.76	3.15 ng/ul	342552	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	95
2		13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	83
3		1-Heneicosyl formate	340	C22H44O2	077899-03-7	81
4		Cyclooctacosane	392	C28H56	000297-24-5	81
5		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\
Data File : BN009510.D
Acq On : 31 Jan 2020 14:37
Operator : CG/JU
Sample : L1271-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GASZ9

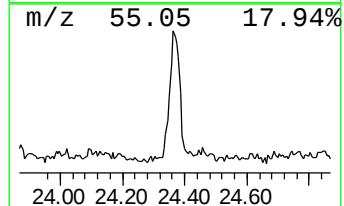
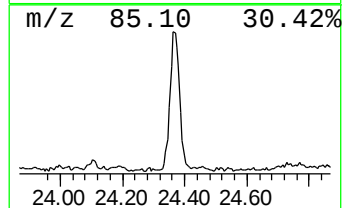
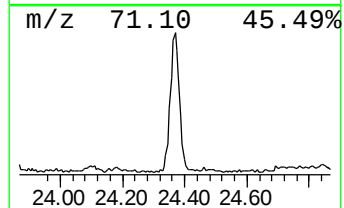
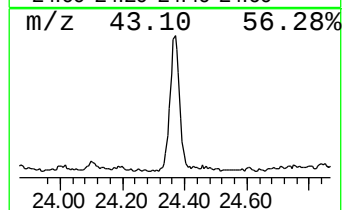
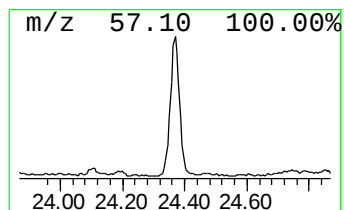
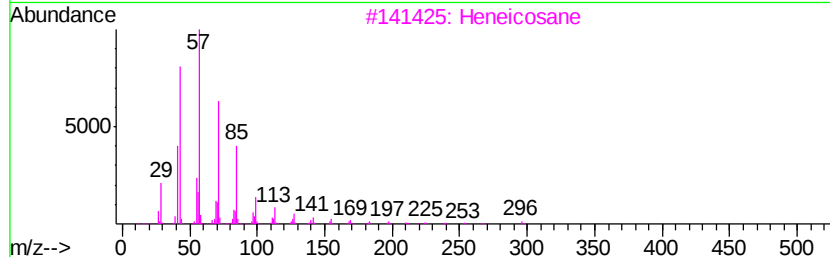
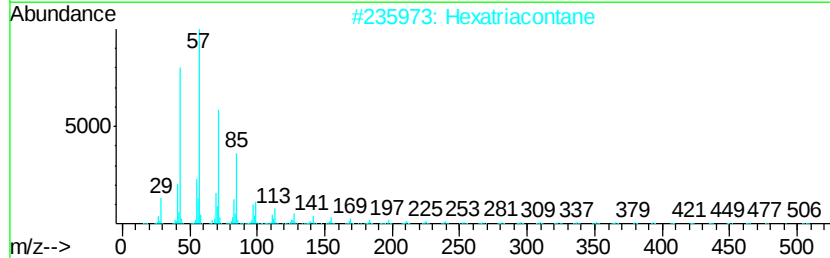
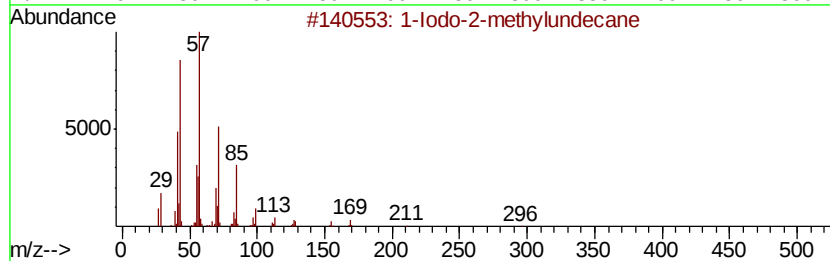
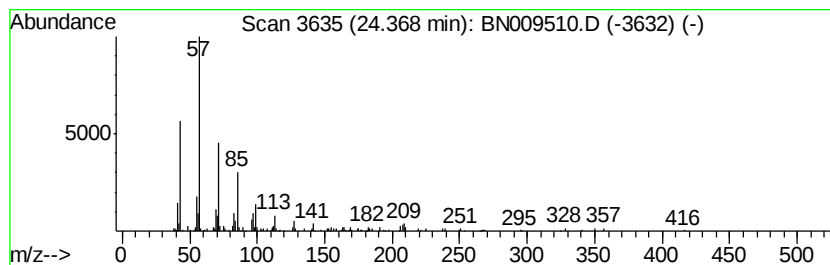
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1-Iodo-2-methylundecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.37	3.73 ng/ul	405509	Perylene-d12	23.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	83
2		Hexatriacontane	507	C36H74	000630-06-8	83
3		Heneicosane	296	C21H44	000629-94-7	83
4		Eicosane	282	C20H42	000112-95-8	74
5		Heneicosane	296	C21H44	000629-94-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN013020\
 Data File : BN009510.D
 Acq On : 31 Jan 2020 14:37
 Operator : CG/JU
 Sample : L1271-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GASZ9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN012220MA.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
n-Hexadecanoic acid	17.77	4.7	ng/ul	501791	4	16.84	2149880	20.0
E-14-Hexadecenal	18.64	2.3	ng/ul	241582	4	16.84	2149880	20.0
(DEL) Alkane: Str...	20.79	10.7	ng/ul	1228890	5	21.05	2299240	20.0
(DEL) Alkane: Str...	21.66	2.8	ng/ul	317926	5	21.05	2299240	20.0
(DEL) Alkane: Str...	22.09	2.7	ng/ul	314503	5	21.05	2299240	20.0
1,5,9-Undecatrien...	22.21	2.6	ng/ul	282787	6	23.16	2174710	20.0
(DEL) Alkane: Str...	22.57	4.2	ng/ul	455243	6	23.16	2174710	20.0
(DEL) Alkane: Str...	23.10	4.7	ng/ul	508458	6	23.16	2174710	20.0
(DEL) Alkane: Str...	23.69	4.6	ng/ul	501623	6	23.16	2174710	20.0
(DEL) Alkane: Cyc...	23.76	3.1	ng/ul	342552	6	23.16	2174710	20.0
1-Iodo-2-methylun...	24.37	3.7	ng/ul	405509	6	23.16	2174710	20.0