

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
 Data File : BN003250.D
 Acq On : 22 Oct 2018 21:33
 Operator : MJ/SJ
 Sample : J5568-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AJ8

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BN101918MA.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	6.805	644	649	667	rBV2	251968	643823	21.92%	1.983%
2	6.952	667	674	685	rVB	216382	363290	12.37%	1.119%
3	7.146	700	707	725	rBV	278392	487415	16.60%	1.501%
4	7.605	779	785	796	rBV	249005	420924	14.33%	1.297%
5	8.328	902	908	929	rBV	324669	645677	21.99%	1.989%
6	8.757	975	981	994	rBV	268101	478750	16.30%	1.475%
7	9.475	1097	1103	1114	rBV	228585	416575	14.19%	1.283%
8	10.022	1190	1196	1218	rBV	298010	689012	23.46%	2.122%
9	10.369	1248	1255	1266	rBV	424033	724835	24.68%	2.233%
10	10.528	1276	1282	1306	rBV	278797	610940	20.80%	1.882%
11	13.657	1808	1814	1819	rBV	750519	1064791	36.26%	3.280%
12	13.710	1819	1823	1836	rVB	179979	286838	9.77%	0.884%
13	13.934	1855	1861	1872	rBV	929376	1393820	47.46%	4.293%
14	14.245	1907	1914	1925	rVB2	688273	1067828	36.36%	3.289%
15	14.522	1955	1961	1986	rBV2	131276	463023	15.77%	1.426%
16	15.245	2066	2084	2100	rBV	1198395	2075455	70.68%	6.393%
17	15.404	2106	2111	2122	rBV	282579	491074	16.72%	1.513%
18	16.416	2279	2283	2290	rBV	61492	93928	3.20%	0.289%
19	16.998	2376	2382	2391	rVB2	959998	1346092	45.84%	4.146%
20	17.098	2393	2399	2410	rVB	1328715	1958723	66.70%	6.034%
21	17.433	2448	2456	2460	rBV6	13419	38742	1.32%	0.119%
22	17.786	2509	2516	2524	rBV2	54276	116993	3.98%	0.360%
23	17.916	2531	2538	2556	rBV	1574442	2419881	82.41%	7.454%
24	18.580	2647	2651	2661	rBV4	31275	58285	1.98%	0.180%
25	18.804	2682	2689	2701	rBV5	57129	202629	6.90%	0.624%
26	19.016	2722	2725	2727	rBV3	29520	40343	1.37%	0.124%
27	19.045	2727	2730	2732	rVV	79081	106510	3.63%	0.328%
28	19.080	2732	2736	2751	rVV3	198171	535935	18.25%	1.651%
29	19.192	2751	2755	2768	rVB2	145579	243397	8.29%	0.750%
30	19.404	2785	2791	2798	rBV2	1650127	2365987	80.57%	7.288%
31	20.051	2897	2901	2906	rBV4	43651	68362	2.33%	0.211%
32	20.174	2919	2922	2924	rBV4	24391	30068	1.02%	0.093%
33	20.451	2965	2969	2972	rBV	48250	56733	1.93%	0.175%
34	20.916	3045	3048	3051	rBV	43779	45137	1.54%	0.139%

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35	20.957	3051	3055	3068	rVV5	93650	274110	9.33%	0.844%
36	21.116	3080	3082	3087	rVB	37854	37498	1.28%	0.116%
37	21.210	3093	3098	3108	rVB	1355159	1815980	61.84%	5.594%
38	21.357	3119	3123	3126	rVB	83764	87902	2.99%	0.271%
39	21.780	3192	3195	3200	rVB	101597	114986	3.92%	0.354%
40	22.233	3268	3272	3277	rBV	144398	187647	6.39%	0.578%
41	22.357	3289	3293	3299	rVB	232327	299678	10.21%	0.923%
42	22.727	3351	3356	3362	rVB	163322	231172	7.87%	0.712%
43	23.274	3442	3449	3458	rVB	1671150	2936570	100.00%	9.046%
44	23.415	3466	3473	3486	rBV	915125	1907029	64.94%	5.874%
45	23.892	3549	3554	3562	rVB	170731	320396	10.91%	0.987%
46	24.645	3670	3682	3702	rBV6	292686	1548606	52.74%	4.770%
47	24.998	3732	3742	3760	rBV6	124693	650648	22.16%	2.004%

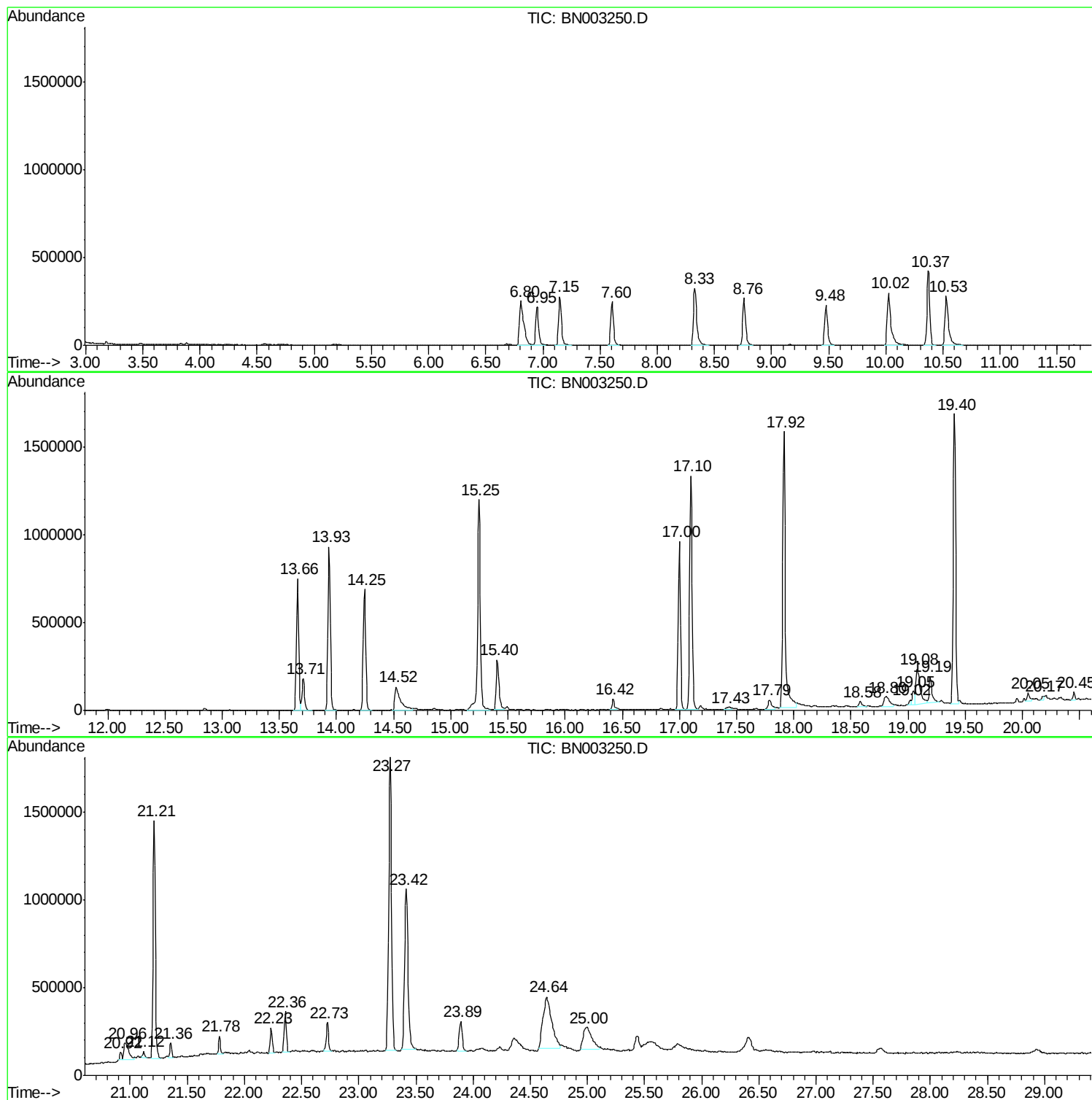
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION

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



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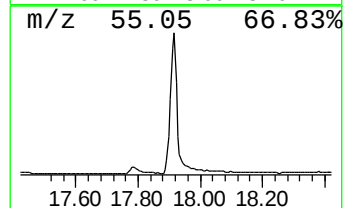
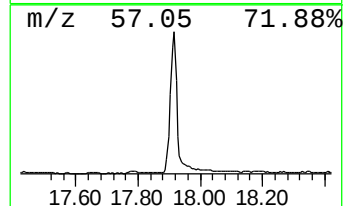
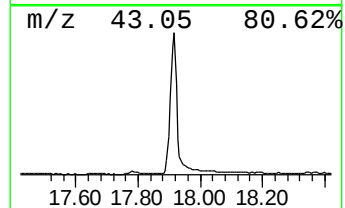
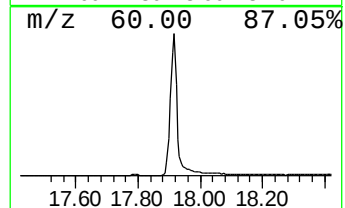
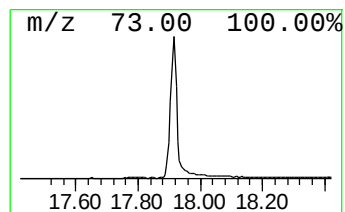
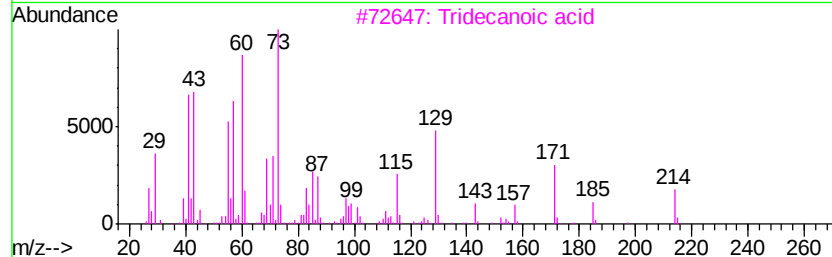
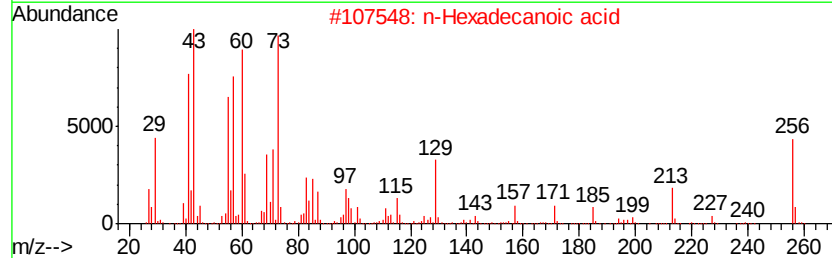
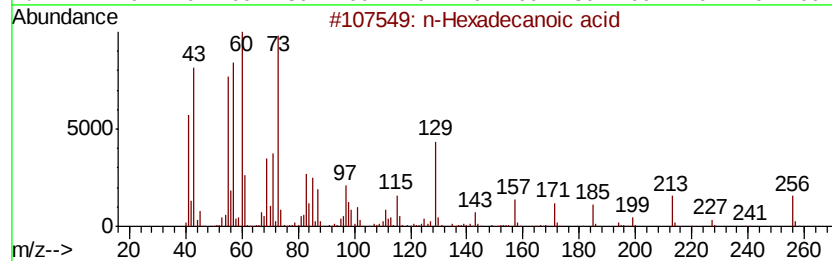
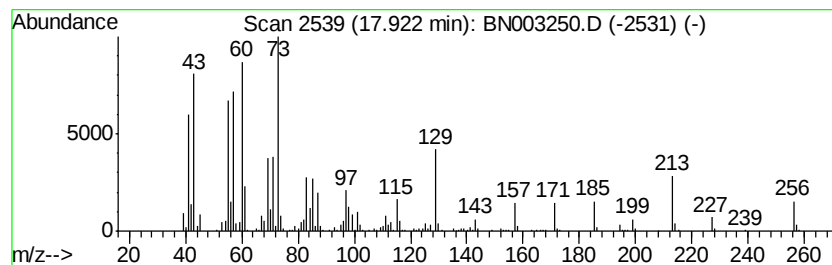
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN101918MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.92	35.95 ng/ul	2419880	Phenanthrene-d10	17.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99	
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98	
3	Tridecanoic acid	214	C13H26O2	000638-53-9	92	
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	91	
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	91	



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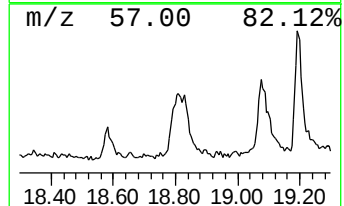
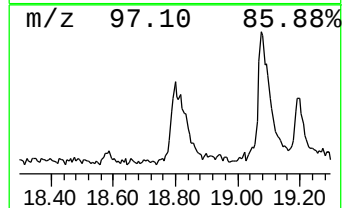
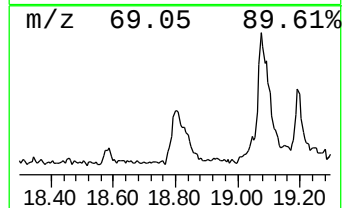
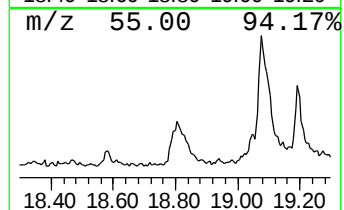
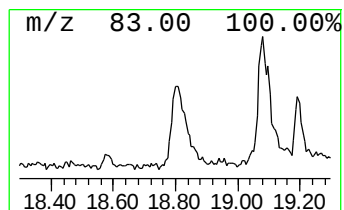
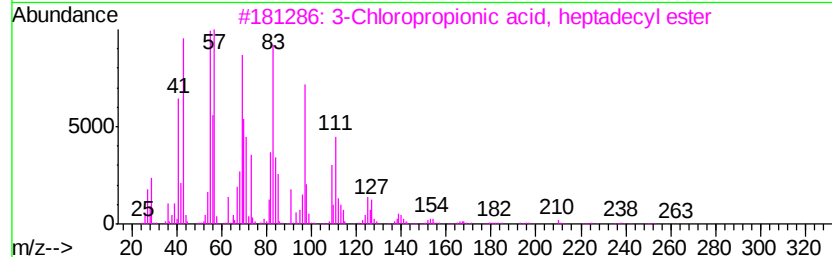
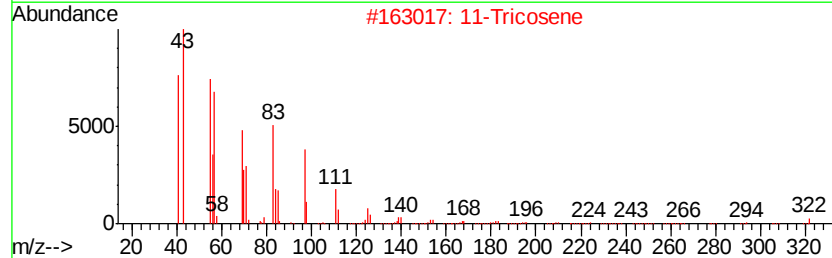
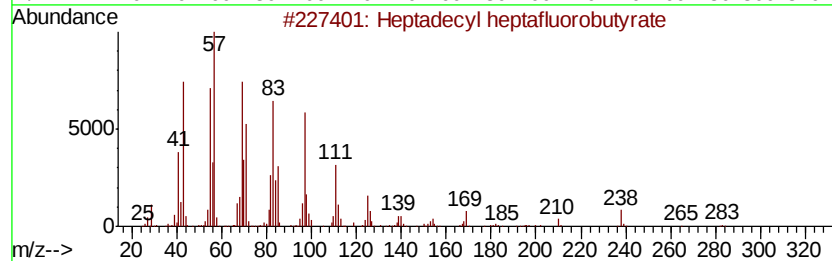
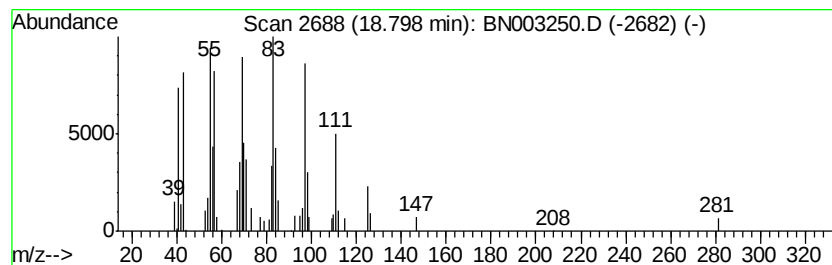
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Peak Number 2 Heptadecyl heptafluorobutyrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.80	3.01 ng/ul	202629	Phenanthrene-d10	17.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl	heptafluorobutyrate	452	C21H35F7O2	959085-66-6	87
2	11-Tricosene		322	C23H46	052078-56-5	87
3	3-Chloropropionic acid, heptadec...		346	C20H39ClO2	1000283-05-1	74
4	9-Eicosene, (E)-		280	C20H40	074685-29-3	74
5	Fumaric acid, 2-chloropropyl tri...		374	C20H35ClO4	1000348-57-1	64



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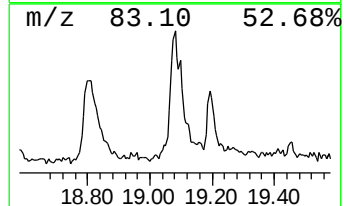
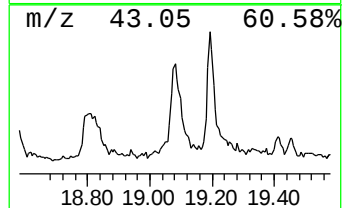
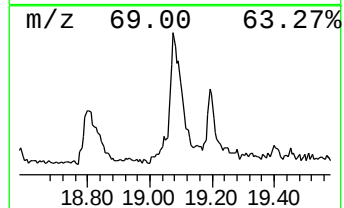
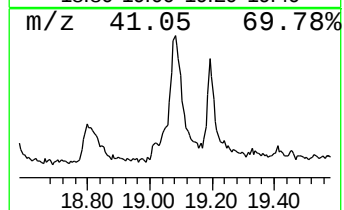
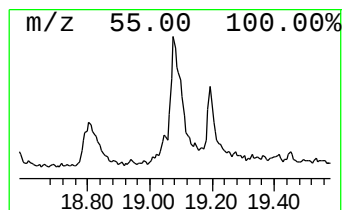
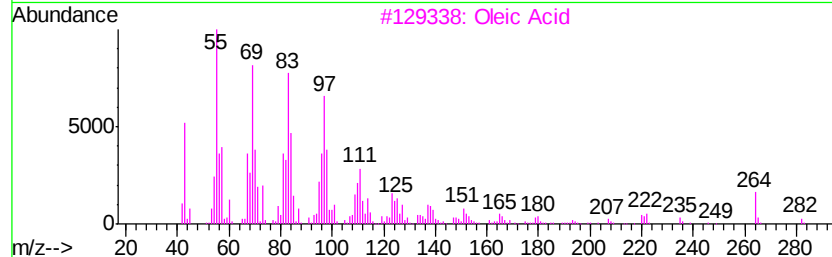
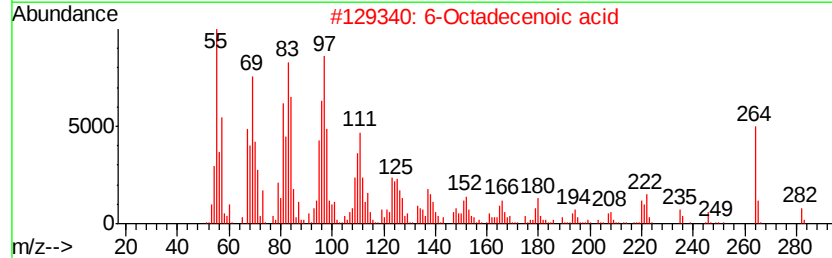
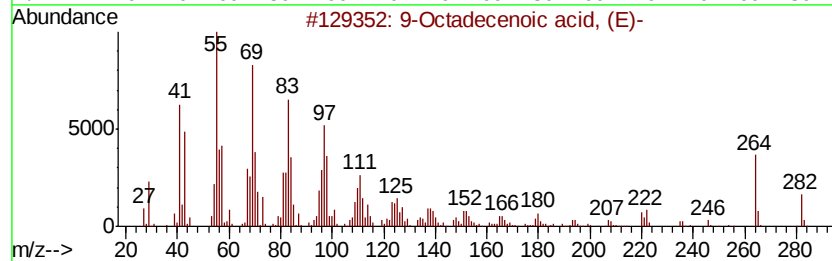
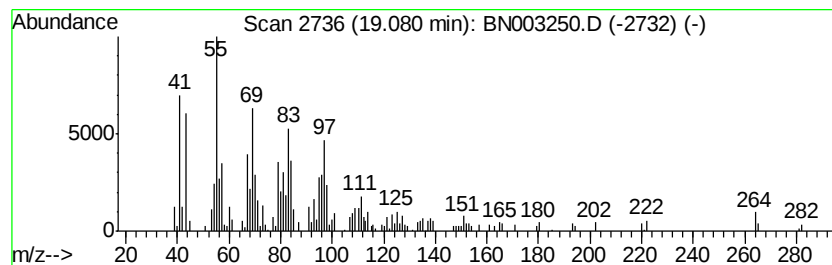
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9-Octadecenoic acid, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.08	7.96 ng/ul	535935	Phenanthrene-d10	17.00

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	96	
2	6-Octadecenoic acid	282	C18H34O2	1000336-66-8	94	
3	Oleic Acid	282	C18H34O2	000112-80-1	93	
4	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	93	
5	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	93	



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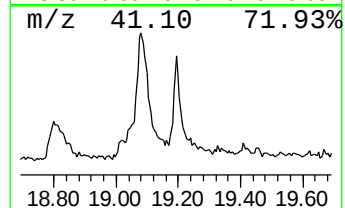
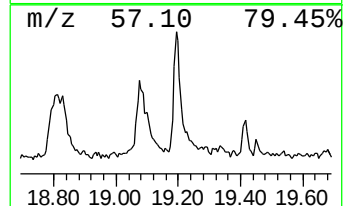
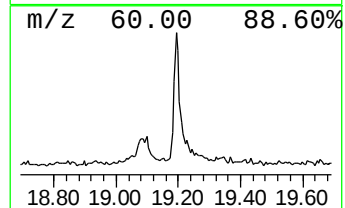
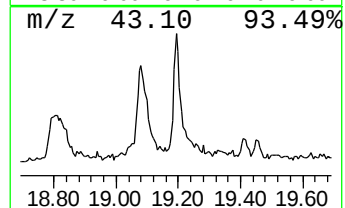
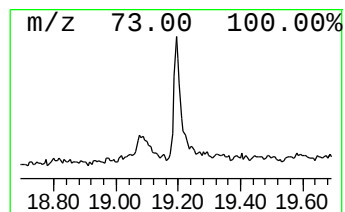
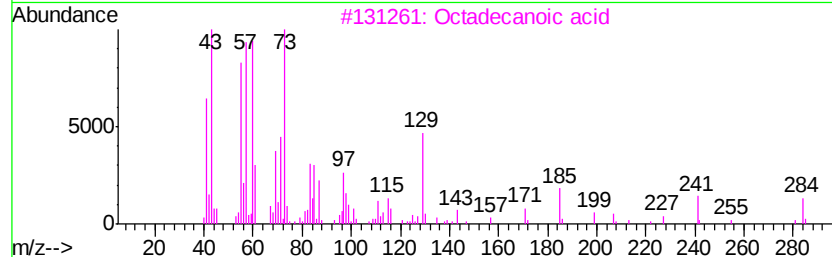
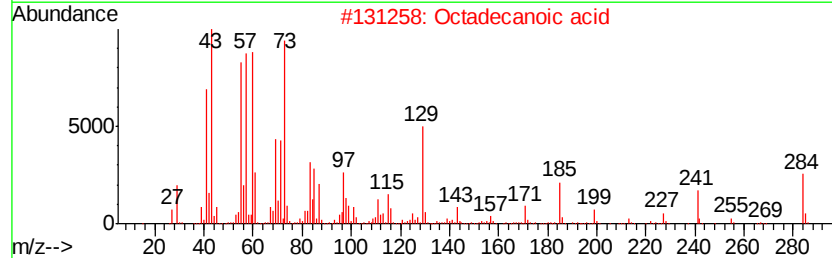
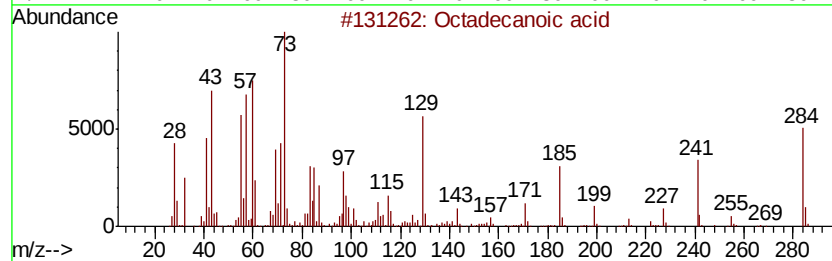
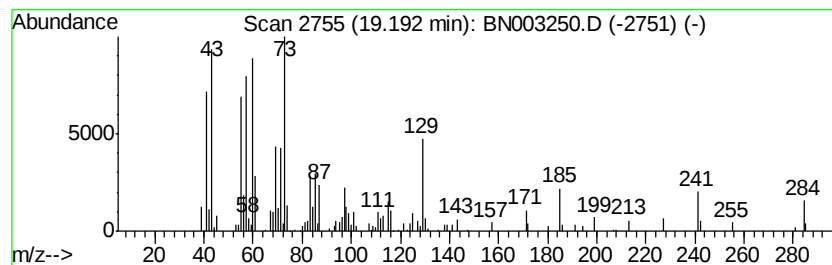
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Peak Number 4 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.19	2.68 ng/ul	243397	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	98
5		Octadecanoic acid	284	C18H36O2	000057-11-4	94



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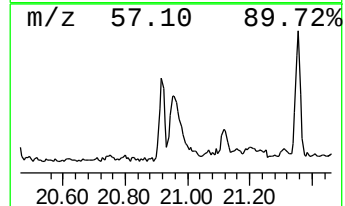
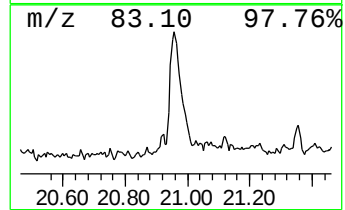
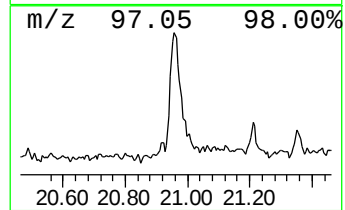
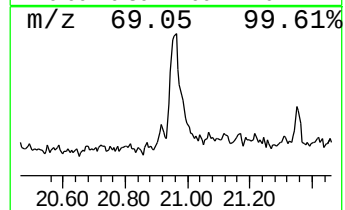
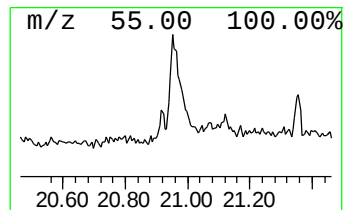
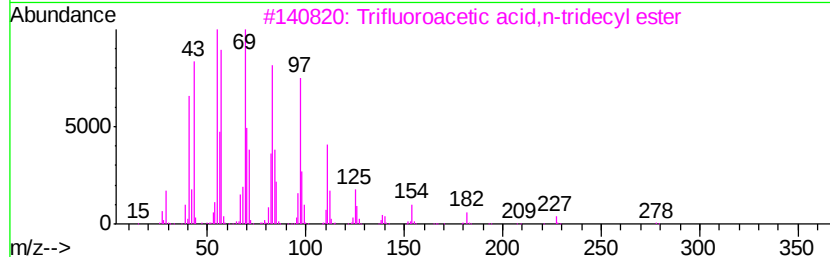
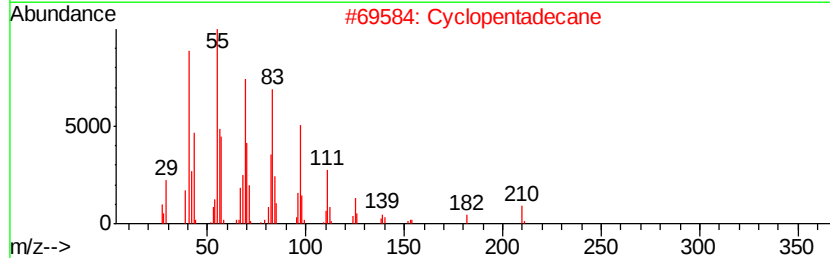
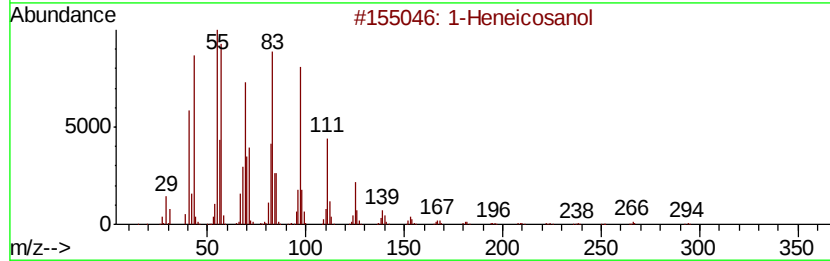
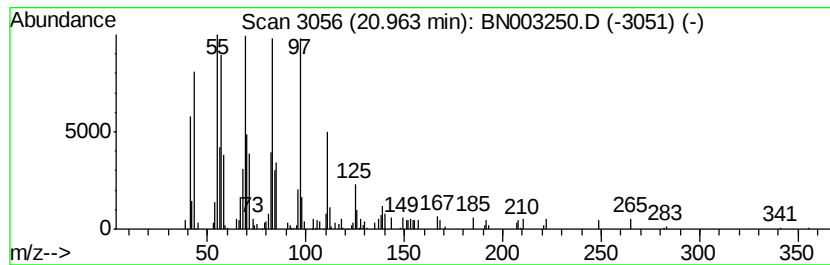
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Heneicosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	3.02 ng/ul	274110	Chrysene-d12	21.21
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Heneicosanol	312 C21H44O	015594-90-8	94
2	Cyclopentadecane	210 C15H30	000295-48-7	91
3	Trifluoroacetic acid,n-tridecyl ...	296 C15H27F3O2	053800-02-5	91
4	1-Octadecanol	270 C18H38O	000112-92-5	90
5	n-Nonadecanol-1	284 C19H40O	001454-84-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003250.D
Acq On : 22 Oct 2018 21:33
Operator : MJ/SJ
Sample : J5568-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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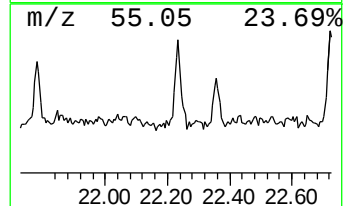
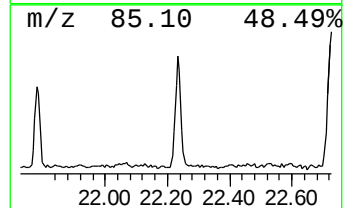
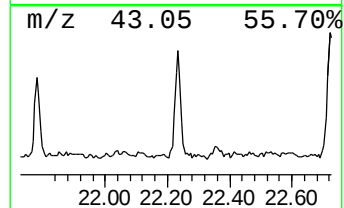
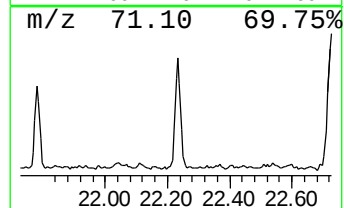
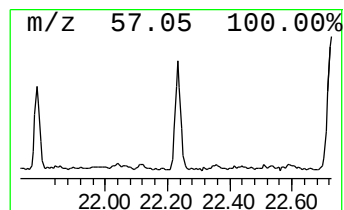
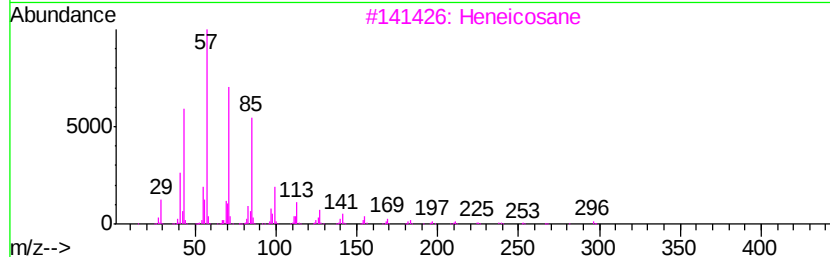
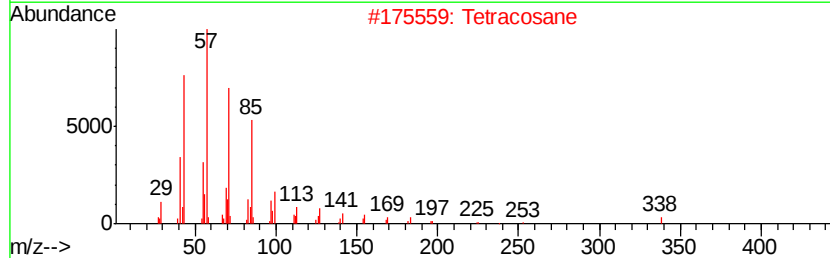
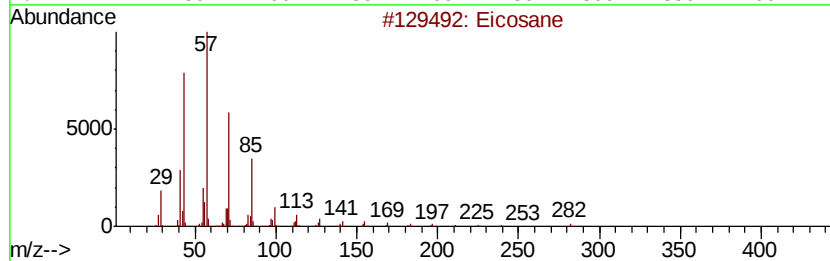
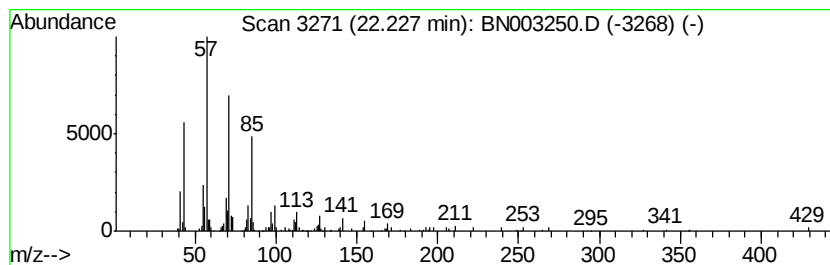
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	2.07 ng/ul	187647	Chrysene-d12	21.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Tetracosane	338	C24H50	000646-31-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5		triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003250.D
Acq On : 22 Oct 2018 21:33
Operator : MJ/SJ
Sample : J5568-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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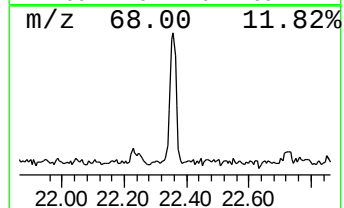
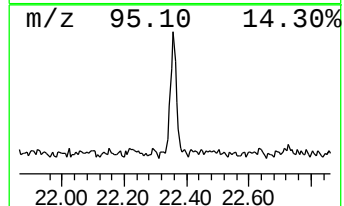
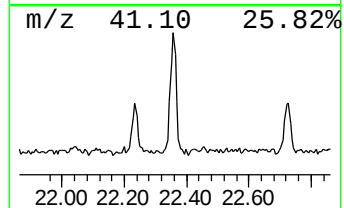
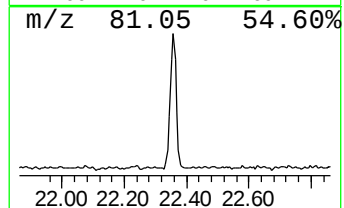
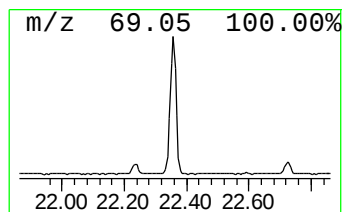
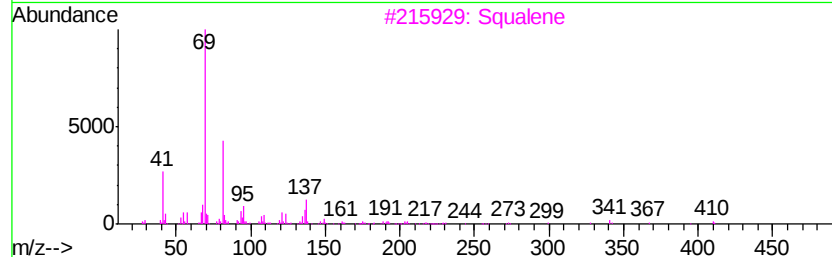
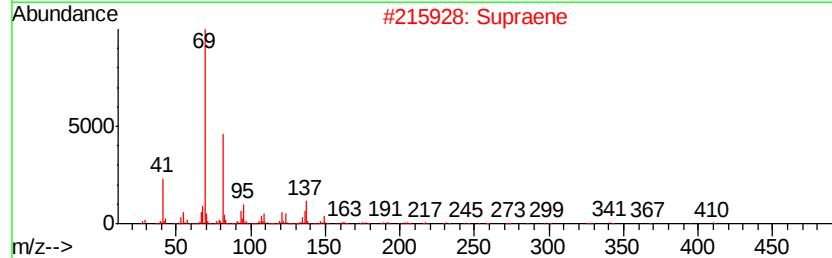
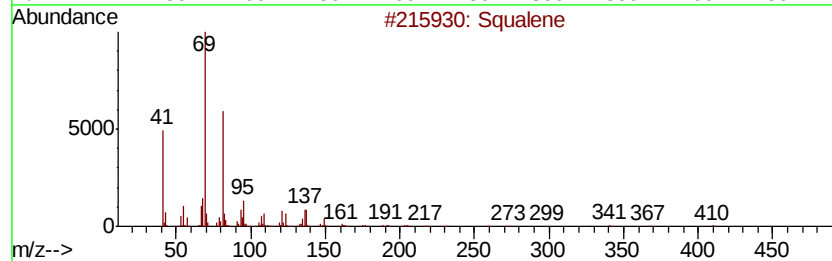
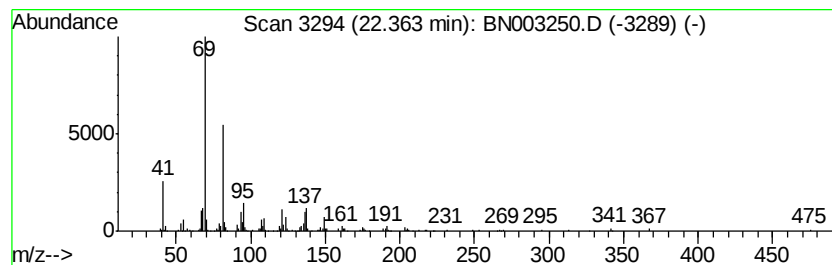
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Squalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	3.14 ng/ul	299678	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	91
2		Supraene	410	C30H50	007683-64-9	91
3		Squalene	410	C30H50	000111-02-4	78
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003250.D
Acq On : 22 Oct 2018 21:33
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Sample : J5568-01
Misc :
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Instrument :
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ClientSampleId :
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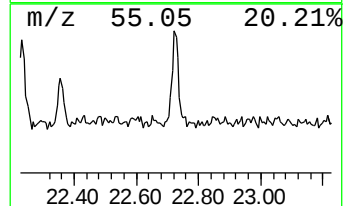
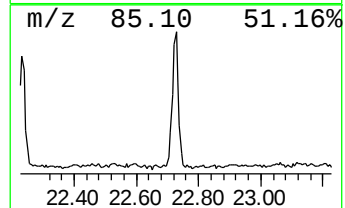
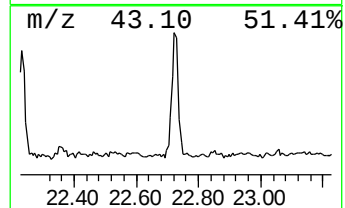
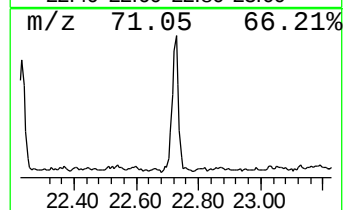
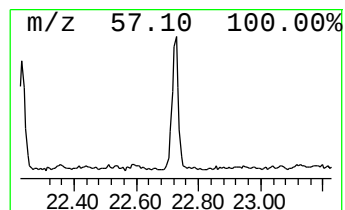
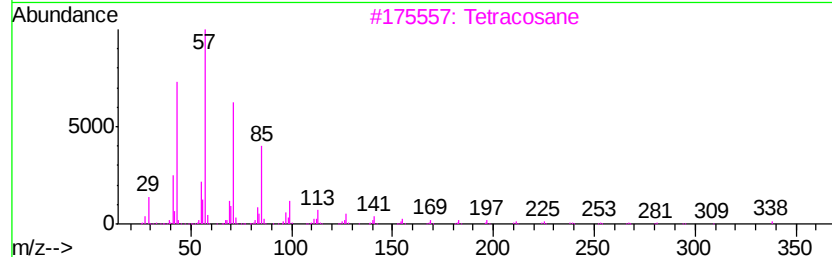
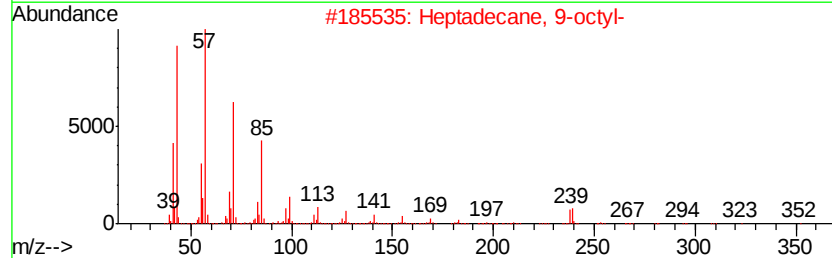
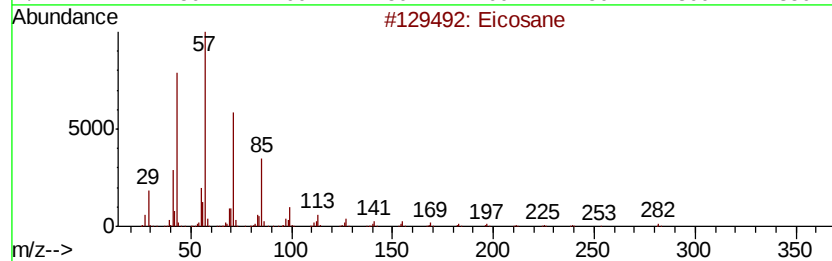
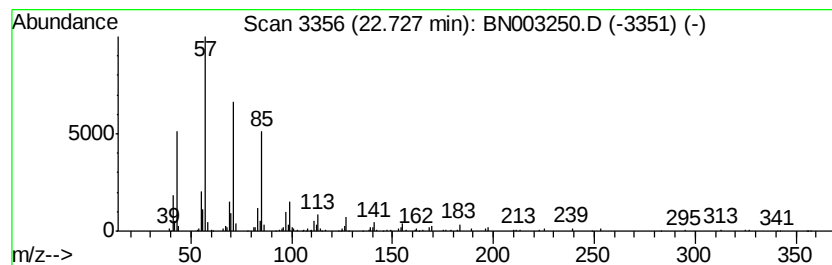
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.73	2.42 ng/ul	231172	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Tetracosane	338	C24H50	000646-31-1	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
Data File : BN003250.D
Acq On : 22 Oct 2018 21:33
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Sample : J5568-01
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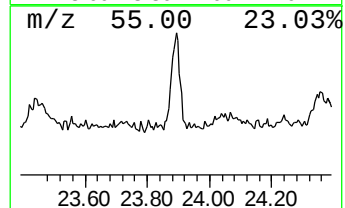
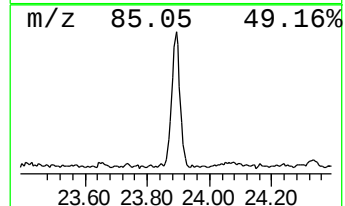
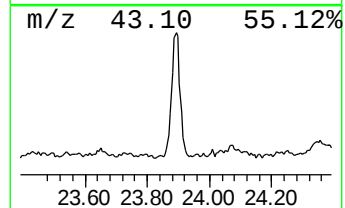
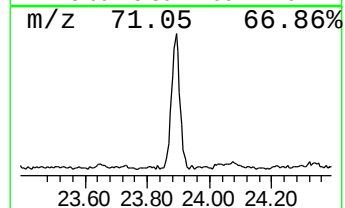
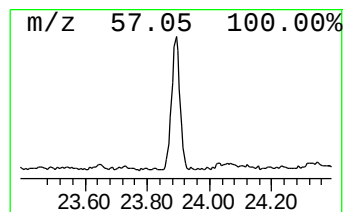
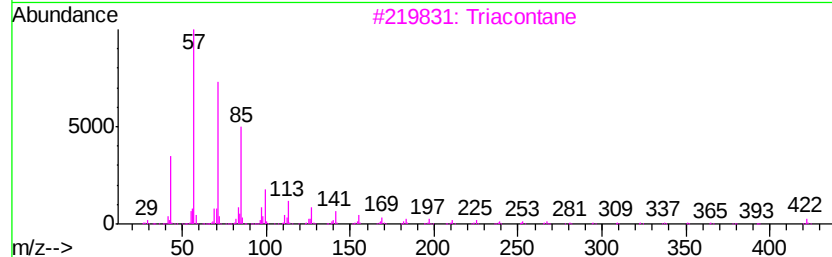
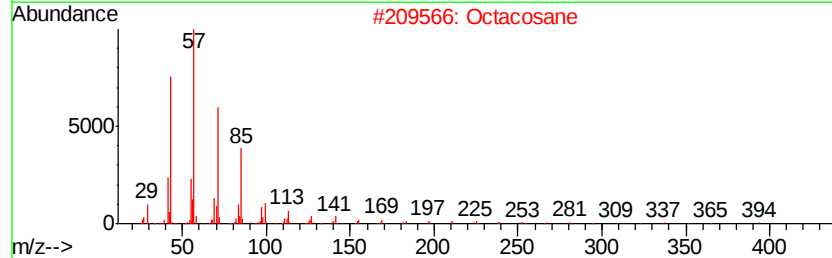
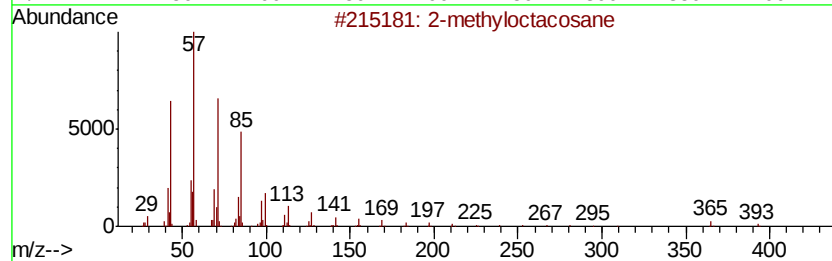
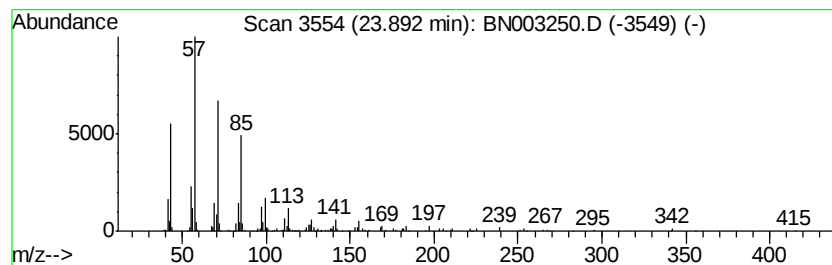
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.89	3.36 ng/ul	320396	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	91
2		Octacosane	394	C28H58	000630-02-4	91
3		Triacotane	422	C30H62	000638-68-6	91
4		Docosane	310	C22H46	000629-97-0	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
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Sample : J5568-01
Misc :
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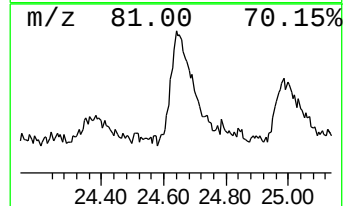
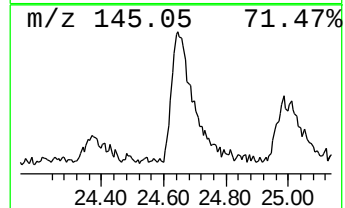
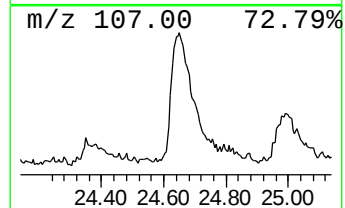
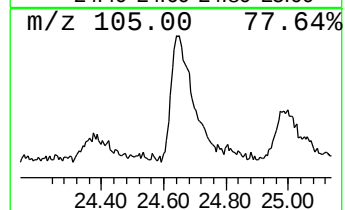
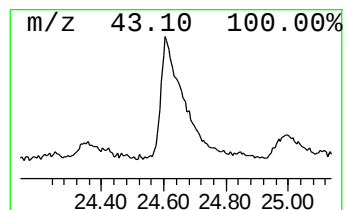
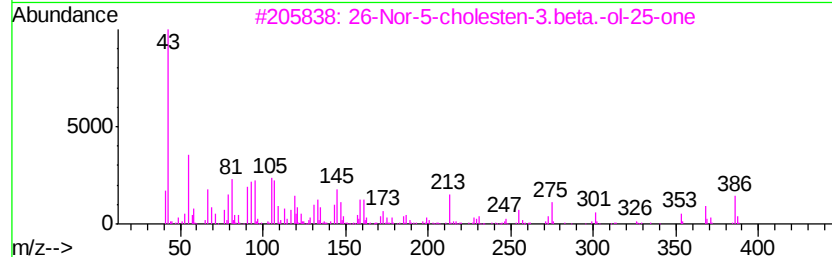
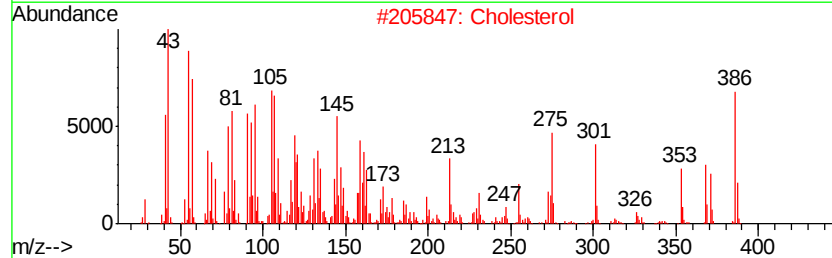
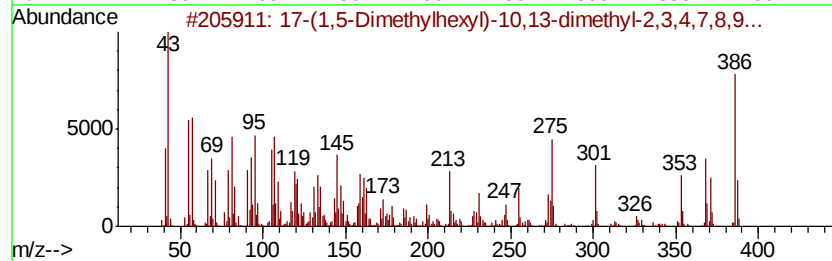
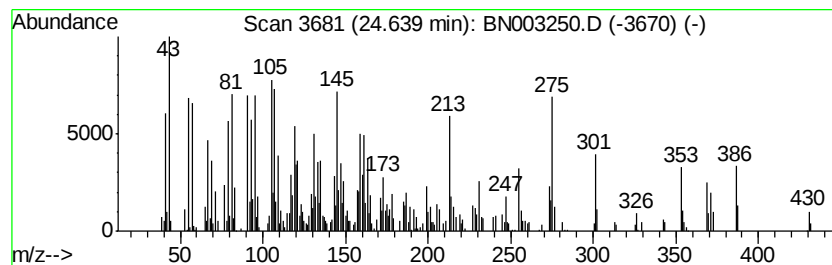
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 17-(1,5-Dimethylhexyl)-10,1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
24.64	16.24 ng/ul	1548610	Perylene-d12	23.42		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-38-4	99
2		Cholesterol	386	C27H46O	000057-88-5	99
3		26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	99
4		Cholesterol	386	C27H46O	000057-88-5	96
5		Cholest-5-en-3-ol, (3.alpha.)-	386	C27H46O	000474-77-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102218\
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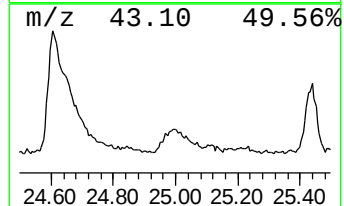
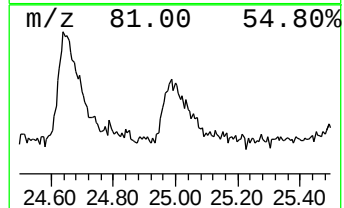
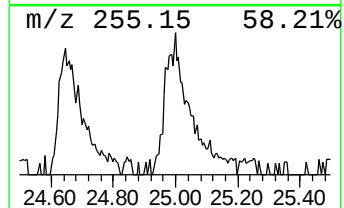
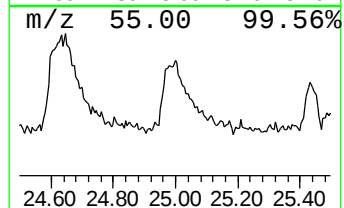
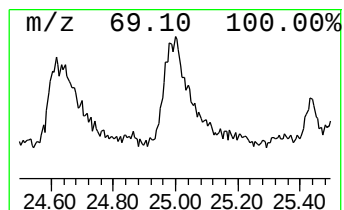
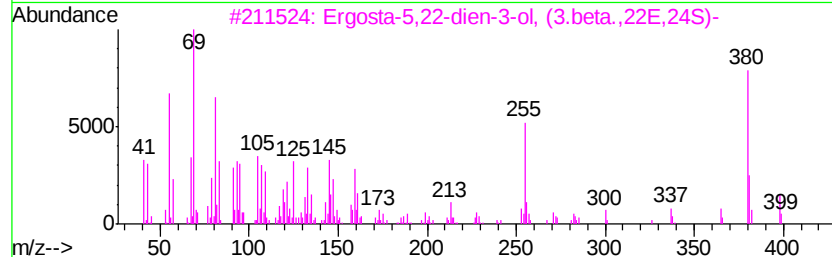
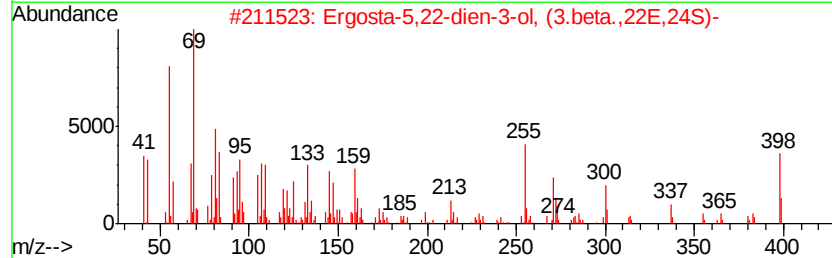
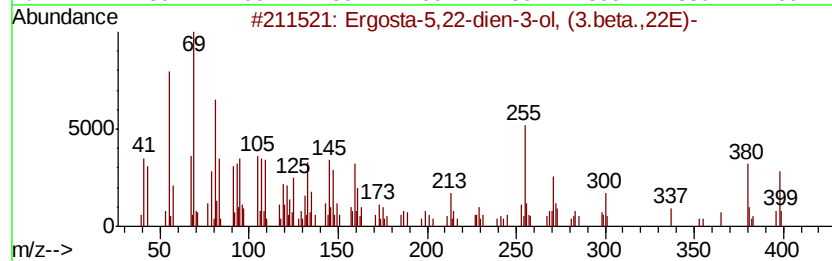
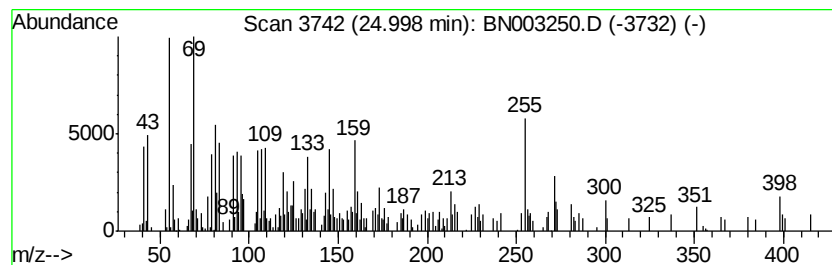
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Ergosta-5,22-dien-3-ol, (3.... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
25.00	6.82 ng/ul	650648	Perylene-d12	23.42		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ergosta-5,22-dien-3-ol, (3.beta....	398	C28H46O	000474-67-9	96	
2	Ergosta-5,22-dien-3-ol, (3.beta....	398	C28H46O	017472-78-5	91	
3	Ergosta-5,22-dien-3-ol, (3.beta....	398	C28H46O	017472-78-5	87	
4	8(14),22-Ergostadienol	398	C28H46O	051502-42-2	70	
5	Spiro[isobenzofuran-1(3H),9'-[9H...	300	C20H12O3	000596-24-7	50	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102218\
Data File : BN003250.D
Acq On : 22 Oct 2018 21:33
Operator : MJ/SJ
Sample : J5568-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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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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Heptadecyl heptaf...	18.80	3.0	ng/ul	202629	4	17.00	1346090	20.0
9-Octadecenoic ac...	19.08	8.0	ng/ul	535935	4	17.00	1346090	20.0
Octadecanoic acid	19.19	2.7	ng/ul	243397	5	21.21	1815980	20.0
1-Heneicosanol	20.96	3.0	ng/ul	274110	5	21.21	1815980	20.0
(DEL) Alkane: Str...	22.23	2.1	ng/ul	187647	5	21.21	1815980	20.0
Squalene	22.36	3.1	ng/ul	299678	6	23.42	1907030	20.0
(DEL) Alkane: Str...	22.73	2.4	ng/ul	231172	6	23.42	1907030	20.0
(DEL) Alkane: Str...	23.89	3.4	ng/ul	320396	6	23.42	1907030	20.0
17-(1,5-Dimethylh...	24.64	16.2	ng/ul	1548610	6	23.42	1907030	20.0
Ergosta-5,22-dien...	25.00	6.8	ng/ul	650648	6	23.42	1907030	20.0