

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
 Data File : BP011507.D
 Acq On : 19 Aug 2022 14:53
 Operator : CG/JU
 Sample : N4210-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGC49

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011507.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.070	11	14	20	rVB	29263	35620	1.48%	0.092%
2	3.487	82	85	99	rBV	195107	307921	12.78%	0.794%
3	3.587	99	102	108	rVB3	20332	28231	1.17%	0.073%
4	3.693	115	120	127	rVB	31745	48976	2.03%	0.126%
5	3.787	131	136	145	rVB	74668	110352	4.58%	0.284%
6	4.011	168	174	179	rBV	32976	48004	1.99%	0.124%
7	4.217	204	209	214	rVB2	13439	20019	0.83%	0.052%
8	4.417	239	243	248	rBV	18330	24931	1.03%	0.064%
9	4.846	311	316	322	rVB	34615	48383	2.01%	0.125%
10	5.581	434	441	447	rVB6	9391	16129	0.67%	0.042%
11	6.746	632	639	652	rVB	404200	659406	27.36%	1.700%
12	6.905	660	666	679	rVB	306699	479029	19.87%	1.235%
13	7.093	691	698	707	rBV	414163	656758	27.25%	1.693%
14	7.181	707	713	715	rBV4	7684	11835	0.49%	0.031%
15	7.558	768	777	785	rBV	474287	726263	30.13%	1.872%
16	8.281	892	900	913	rBV	434870	726483	30.14%	1.873%
17	8.393	915	919	927	rVB2	10227	16732	0.69%	0.043%
18	8.722	968	975	981	rBV	392951	621512	25.79%	1.602%
19	9.122	1036	1043	1049	rBV	9157	15369	0.64%	0.040%
20	9.440	1089	1097	1110	rBV	327275	542720	22.52%	1.399%
21	9.975	1177	1188	1199	rBV	456079	772668	32.06%	1.992%
22	10.346	1242	1251	1257	rBV	611121	1007164	41.79%	2.596%
23	10.393	1257	1259	1265	rVV	15235	20953	0.87%	0.054%
24	10.499	1269	1277	1295	rVV	330432	613345	25.45%	1.581%
25	11.340	1415	1420	1422	rBV5	10276	14299	0.59%	0.037%
26	11.946	1518	1523	1530	rVV3	9324	18429	0.76%	0.048%
27	12.022	1530	1536	1542	rVB	16190	27171	1.13%	0.070%
28	12.169	1556	1561	1568	rBV5	8208	19548	0.81%	0.050%
29	12.240	1568	1573	1580	rVB3	15928	30025	1.25%	0.077%
30	13.051	1707	1711	1717	rVV4	14515	23940	0.99%	0.062%
31	13.134	1721	1725	1728	rVV6	6586	12007	0.50%	0.031%
32	13.545	1788	1795	1801	rBV3	12266	25466	1.06%	0.066%
33	13.604	1801	1805	1808	rVV3	9403	14442	0.60%	0.037%
34	13.657	1808	1814	1820	rVV	806037	1122547	46.57%	2.894%
35	13.716	1821	1824	1829	rVV5	10907	17889	0.74%	0.046%
36	13.781	1829	1835	1839	rVV5	7937	18321	0.76%	0.047%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
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37	13.922	1851	1859	1874	rBV	919860	1380116	57.26%	3.558%
38	14.140	1892	1896	1900	rVB2	13647	18719	0.78%	0.048%
39	14.228	1904	1911	1919	rVV	892126	1310434	54.37%	3.378%
40	14.292	1919	1922	1931	rVB5	9142	16478	0.68%	0.042%
41	14.463	1945	1951	1971	rBV	286684	530543	22.01%	1.368%
42	14.634	1975	1980	1984	rVV2	13992	26072	1.08%	0.067%
43	14.675	1984	1987	1991	rVB5	13161	16446	0.68%	0.042%
44	14.851	2012	2017	2022	rVB8	5755	12966	0.54%	0.033%
45	14.904	2022	2026	2033	rVB9	10022	20441	0.85%	0.053%
46	15.022	2041	2046	2049	rBV7	5842	12827	0.53%	0.033%
47	15.104	2056	2060	2062	rBV3	12803	18091	0.75%	0.047%
48	15.134	2062	2065	2069	rVB	23547	29431	1.22%	0.076%
49	15.234	2075	2082	2088	rBV	1193136	1624156	67.38%	4.187%
50	15.281	2088	2090	2096	rVB4	15723	21942	0.91%	0.057%
51	15.375	2100	2106	2115	rBV	279270	393650	16.33%	1.015%
52	15.510	2126	2129	2134	rVB7	9176	14976	0.62%	0.039%
53	15.587	2139	2142	2146	rBV2	8911	12470	0.52%	0.032%
54	15.975	2203	2208	2210	rBV3	26548	40169	1.67%	0.104%
55	16.151	2231	2238	2244	rVV10	11969	32392	1.34%	0.084%
56	16.416	2279	2283	2290	rBV10	12357	19889	0.83%	0.051%
57	16.486	2291	2295	2299	rVV3	9184	13197	0.55%	0.034%
58	16.545	2299	2305	2309	rVV9	10296	24801	1.03%	0.064%
59	16.663	2320	2325	2332	rVB5	15261	33280	1.38%	0.086%
60	16.734	2333	2337	2341	rBV6	8082	14288	0.59%	0.037%
61	16.775	2341	2344	2349	rVV4	25785	35817	1.49%	0.092%
62	16.822	2349	2352	2356	rVB3	14579	20230	0.84%	0.052%
63	16.922	2365	2369	2376	rBV2	50911	71460	2.96%	0.184%
64	17.004	2376	2383	2387	rVV	1095770	1546871	64.18%	3.988%
65	17.045	2387	2390	2394	rVV	160499	212294	8.81%	0.547%
66	17.104	2394	2400	2411	rVV	1243147	1826695	75.79%	4.709%
67	17.198	2412	2416	2422	rVV8	13909	32146	1.33%	0.083%
68	17.422	2451	2454	2463	rVB3	14626	24043	1.00%	0.062%
69	17.522	2468	2471	2475	rBV2	20461	32291	1.34%	0.083%
70	17.704	2498	2502	2506	rBV	45417	52342	2.17%	0.135%
71	17.810	2514	2520	2523	rBV3	32386	51295	2.13%	0.132%
72	17.869	2525	2530	2536	rBV2	212289	323768	13.43%	0.835%
73	17.928	2536	2540	2546	rBV	473140	594689	24.67%	1.533%
74	18.069	2560	2564	2569	rBV2	31597	46200	1.92%	0.119%
75	18.210	2584	2588	2595	rBV4	40726	82397	3.42%	0.212%
76	18.380	2614	2617	2620	rBV	24505	27628	1.15%	0.071%

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77	18.422	2621	2624	2628	rVB2	25645	30112	1.25%	0.078%
78	18.786	2682	2686	2689	rBV2	17915	30119	1.25%	0.078%
79	18.828	2689	2693	2699	rVV3	38243	56454	2.34%	0.146%
80	18.880	2699	2702	2705	rVV4	17563	19366	0.80%	0.050%
81	18.922	2706	2709	2717	rVV4	43328	92937	3.86%	0.240%
82	19.039	2724	2729	2733	rBV	446651	575294	23.87%	1.483%
83	19.080	2733	2736	2743	rVB3	62537	81818	3.39%	0.211%
84	19.175	2749	2752	2758	rBV3	82797	125798	5.22%	0.324%
85	19.369	2779	2785	2788	rBV	1657453	2248288	93.28%	5.796%
86	19.916	2876	2878	2881	rVB	58682	46814	1.94%	0.121%
87	20.404	2958	2961	2964	rBV	74025	78260	3.25%	0.202%
88	20.869	3036	3040	3046	rVB4	347839	526896	21.86%	1.358%
89	21.063	3069	3073	3078	rBV	1280047	1332233	55.27%	3.435%
90	21.133	3079	3085	3089	rVV2	1402434	2015211	83.61%	5.195%
91	21.169	3089	3091	3102	rVB	225878	308650	12.81%	0.796%
92	21.739	3185	3188	3191	rBV2	237621	277079	11.50%	0.714%
93	22.004	3230	3233	3238	rBV	361676	431377	17.90%	1.112%
94	22.745	3354	3359	3364	rBV2	494336	922348	38.27%	2.378%
95	22.810	3366	3370	3382	rVB	811857	1805510	74.91%	4.655%
96	23.280	3445	3450	3455	rBV2	1054525	1835974	76.17%	4.733%
97	23.433	3470	3476	3482	rBV2	852666	1607203	66.68%	4.143%
98	24.033	3574	3578	3588	rVB	364337	738023	30.62%	1.903%
99	25.780	3870	3875	3892	rVB5	540043	1672587	69.39%	4.312%
100	26.680	4021	4028	4047	rVB2	620830	2410272	100.00%	6.214%

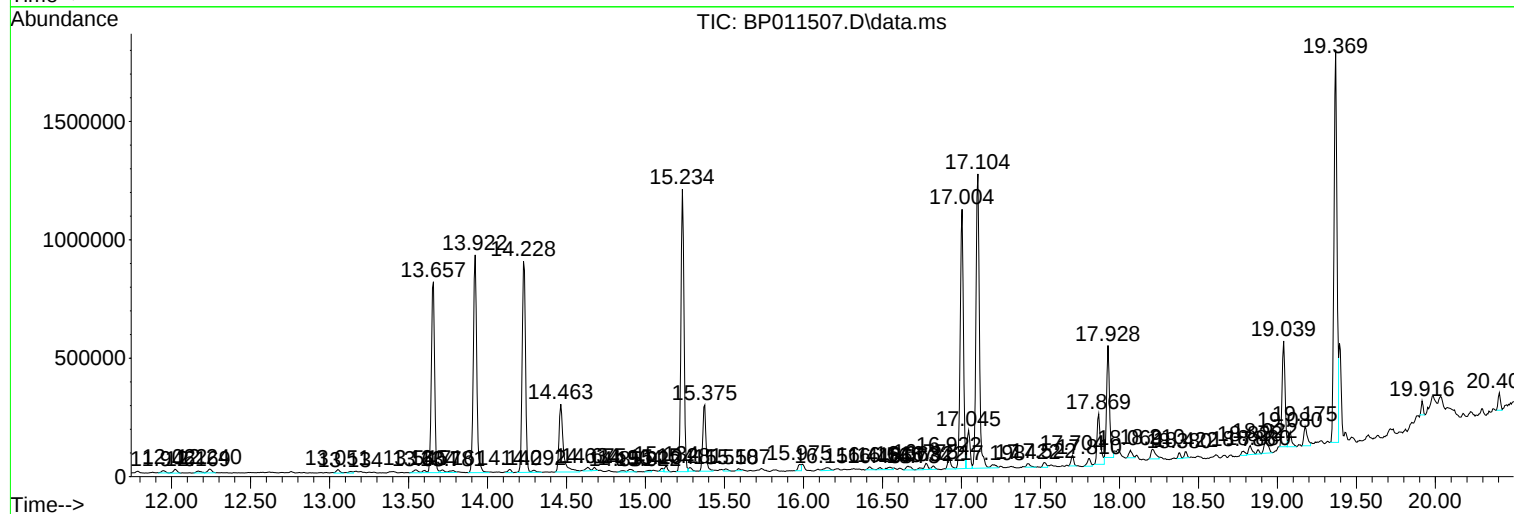
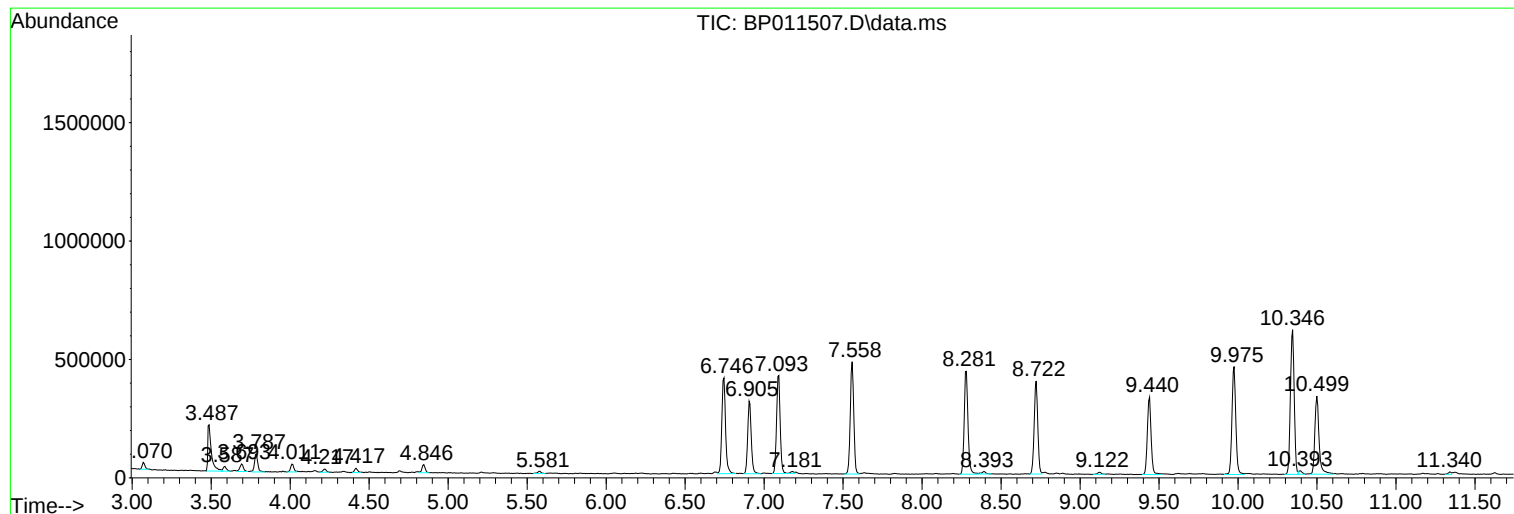
Sum of corrected areas: 38789450

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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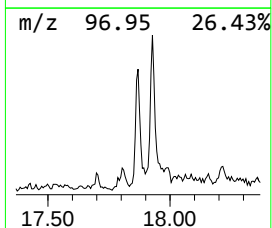
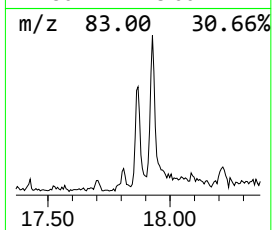
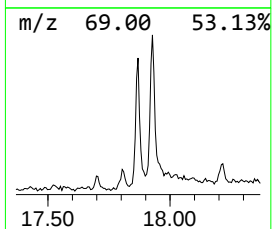
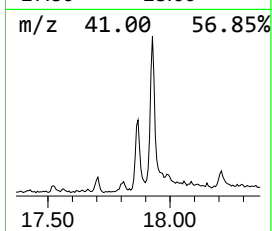
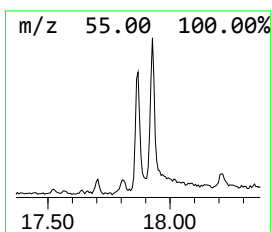
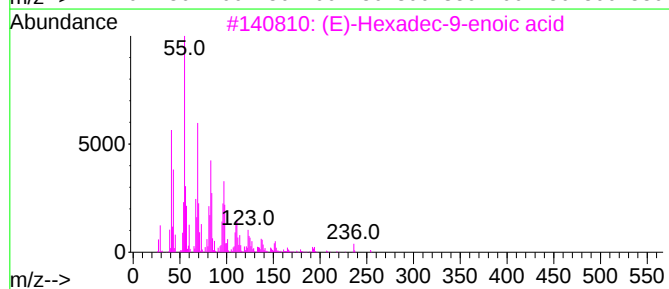
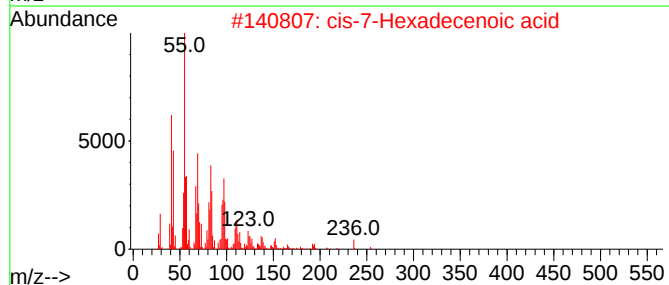
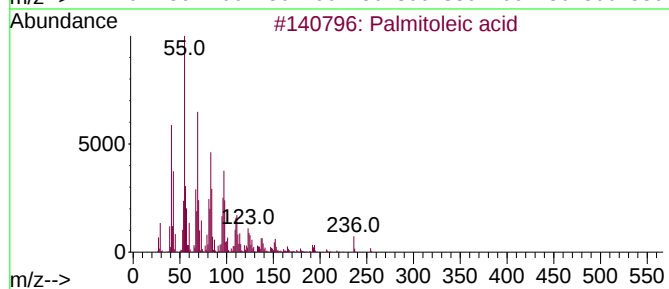
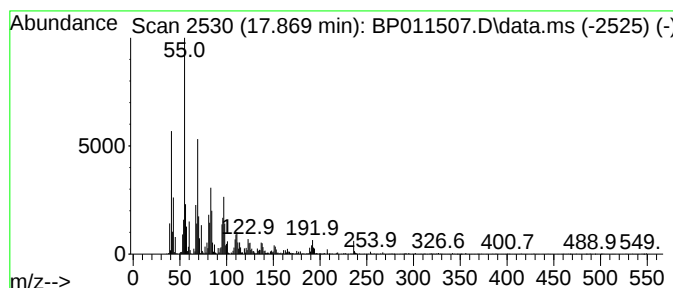
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Palmitoleic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.869	4.19 ng/ul	323768	Phenanthrene-d10	17.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	99
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	99
3			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
4			cis-11-Eicosenoic acid	310	C20H38O2	005561-99-9	99
5			Palmitoleic acid	254	C16H30O2	000373-49-9	98



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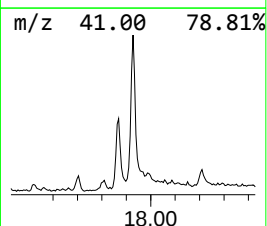
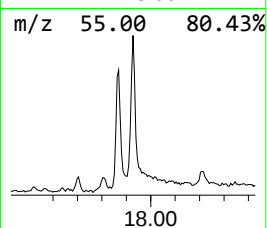
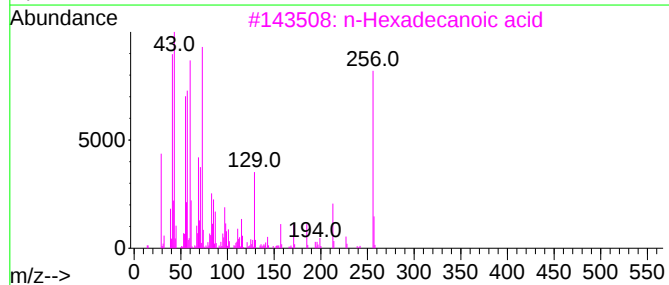
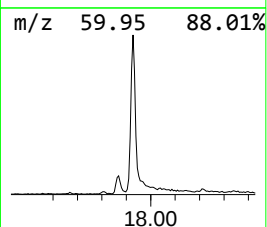
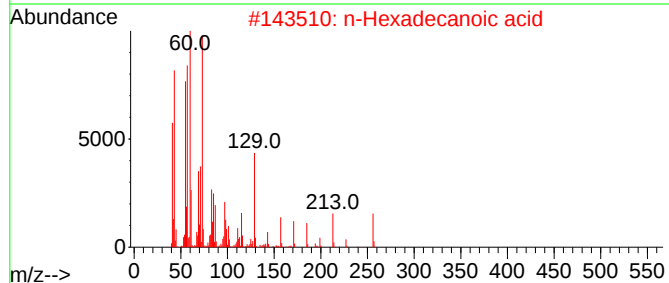
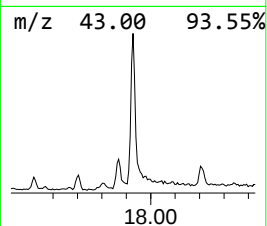
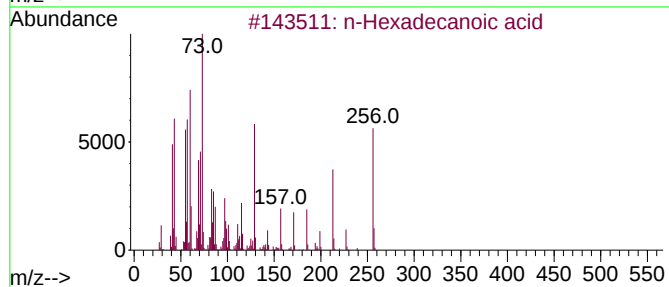
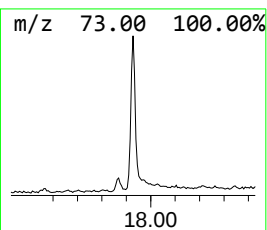
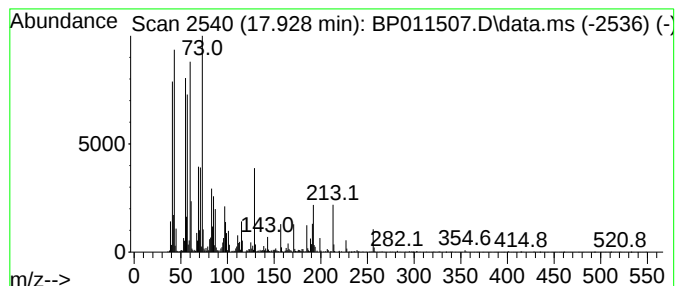
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.928	7.69 ng/ul	594689	Phenanthrene-d10	17.004

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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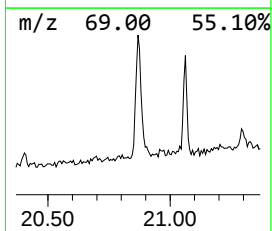
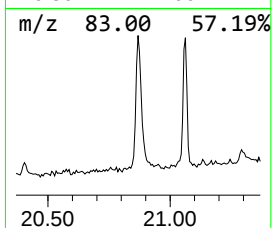
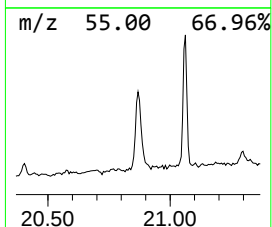
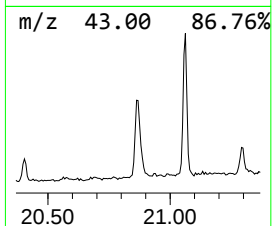
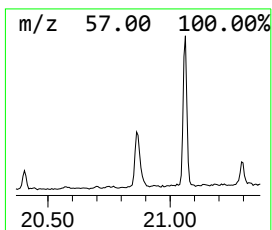
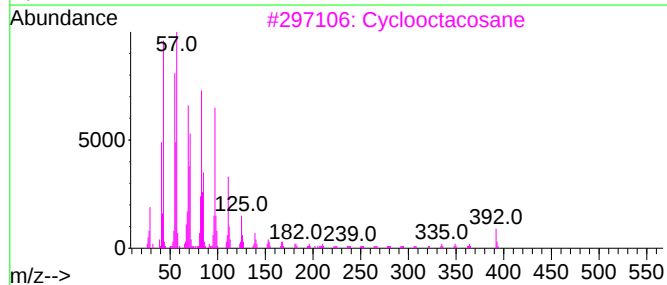
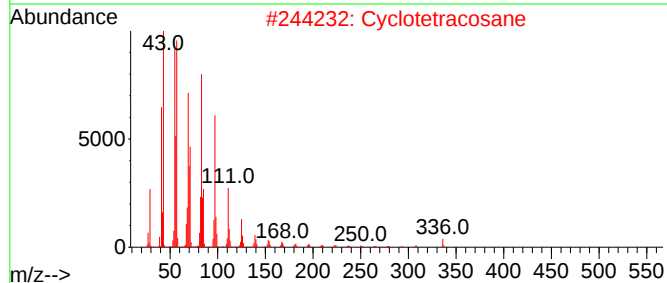
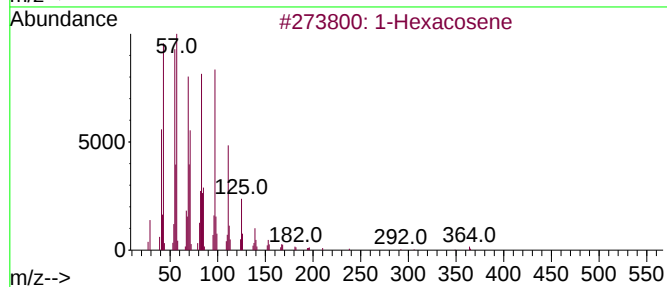
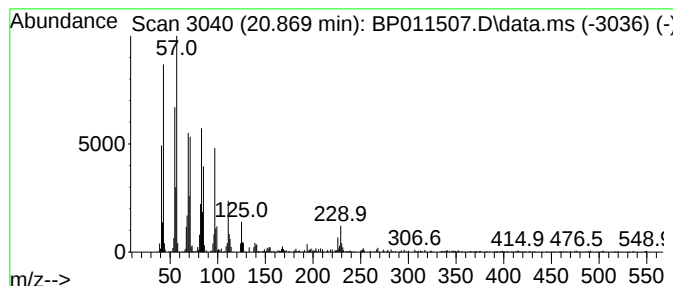
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.869	5.23 ng/ul	526896	Chrysene-d12	21.133	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	98
2	Cyclotetracosane	336	C24H48	000297-03-0	95
3	Cyclooctacosane	392	C28H56	000297-24-5	93
4	Hexadecane, 1-(ethenyloxy)-	268	C18H36O	000822-28-6	91
5	1-Docosene	308	C22H44	001599-67-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011507.D
Acq On : 19 Aug 2022 14:53
Operator : CG/JU
Sample : N4210-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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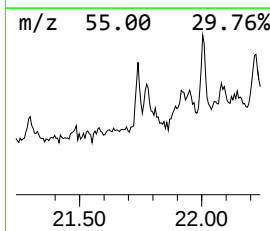
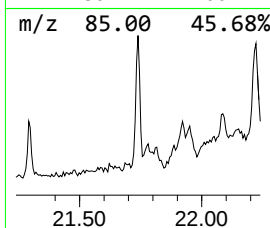
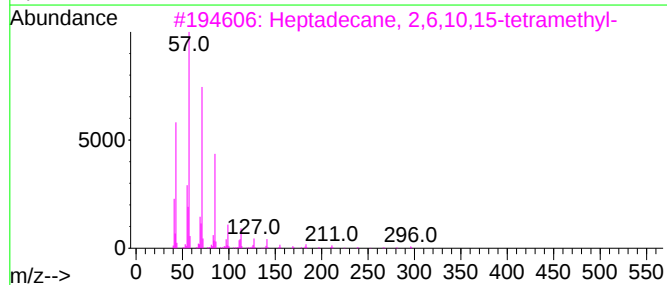
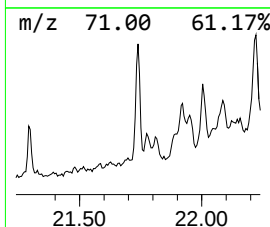
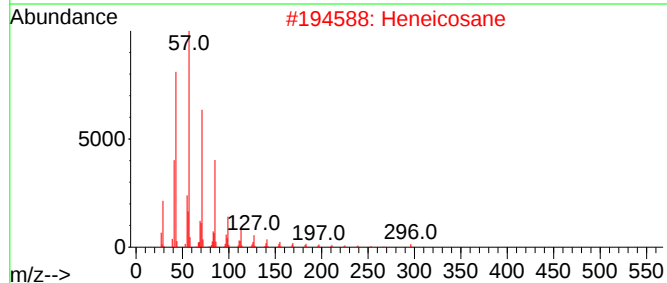
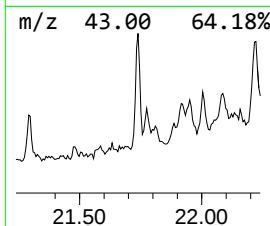
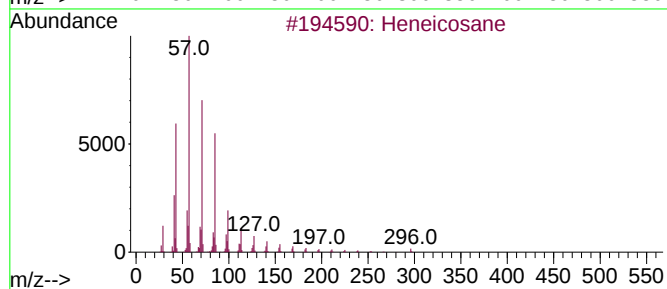
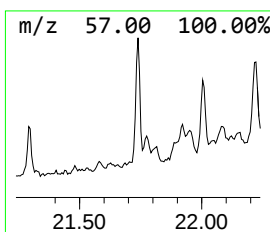
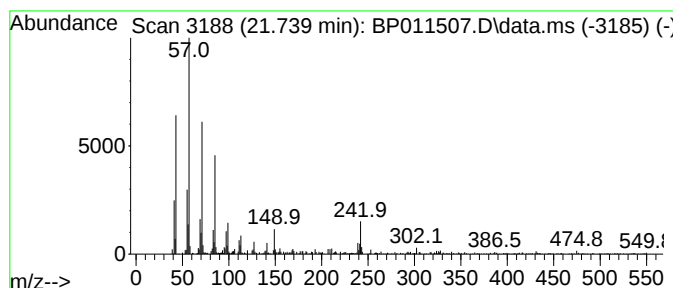
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.739	2.75 ng/ul	277079	Chrysene-d12	21.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			Heneicosane	296	C21H44	000629-94-7	95
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	95
4			Heptadecane	240	C17H36	000629-78-7	95
5			Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011507.D
Acq On : 19 Aug 2022 14:53
Operator : CG/JU
Sample : N4210-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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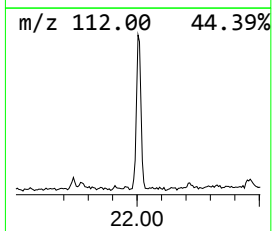
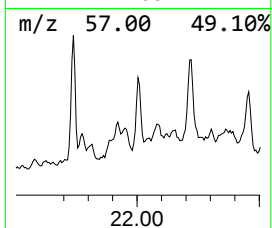
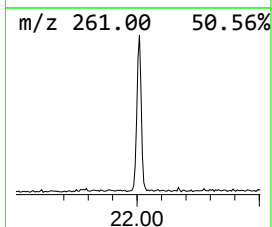
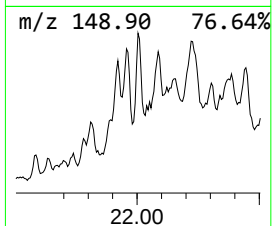
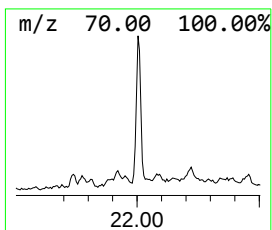
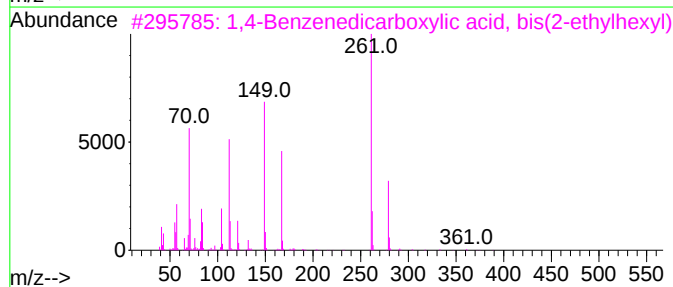
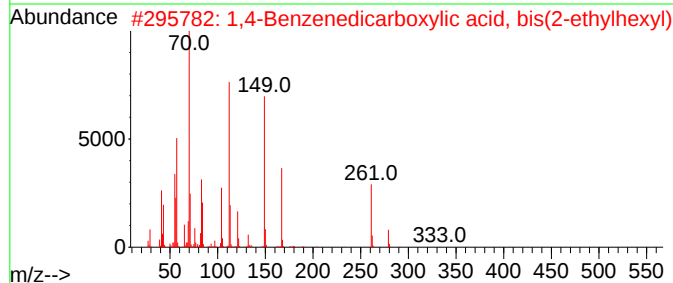
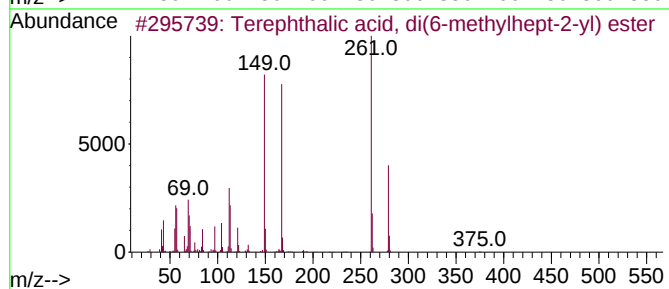
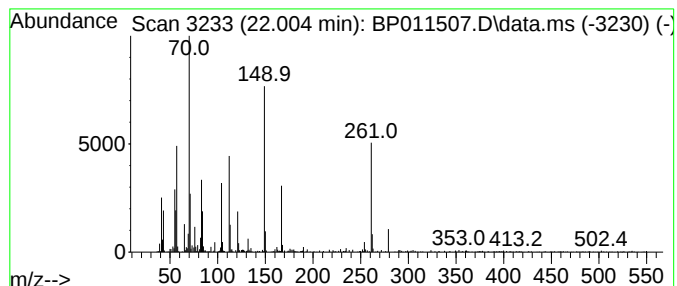
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Terephthalic acid, di(6-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.004	4.28 ng/ul	431377	Chrysene-d12	21.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Terephthalic acid, di(6-methylhe...	390	C24H38O4	1000416-00-0	58
2			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	52
3			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	52
4			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	50
5			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011507.D
Acq On : 19 Aug 2022 14:53
Operator : CG/JU
Sample : N4210-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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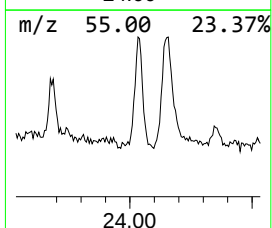
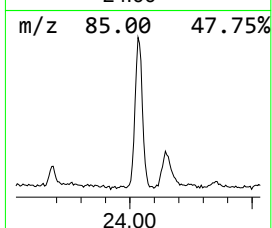
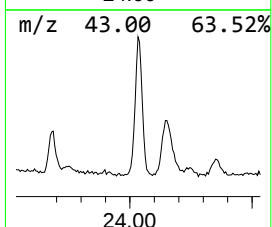
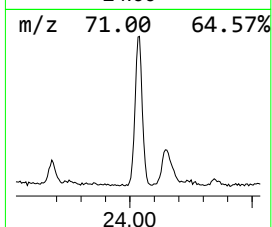
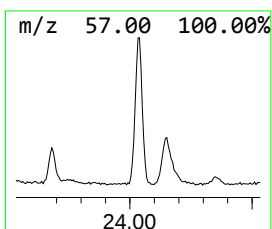
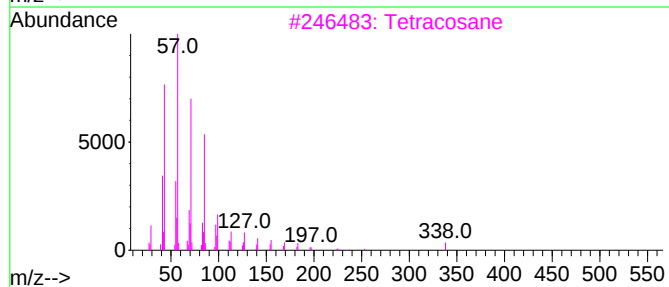
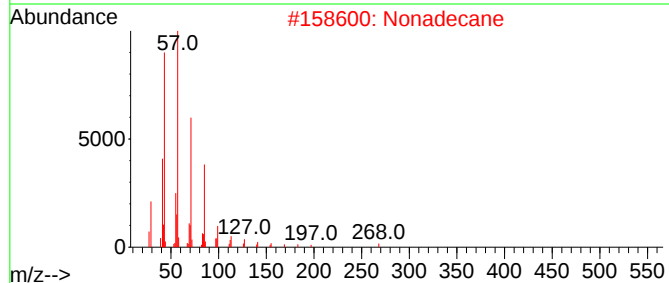
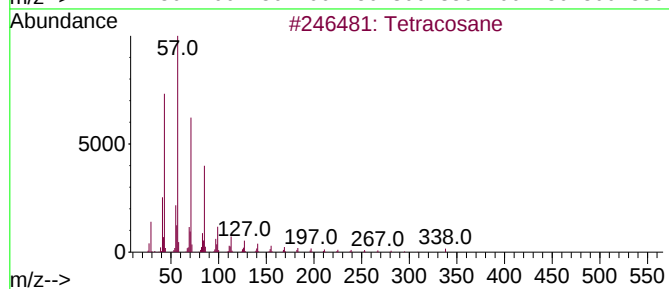
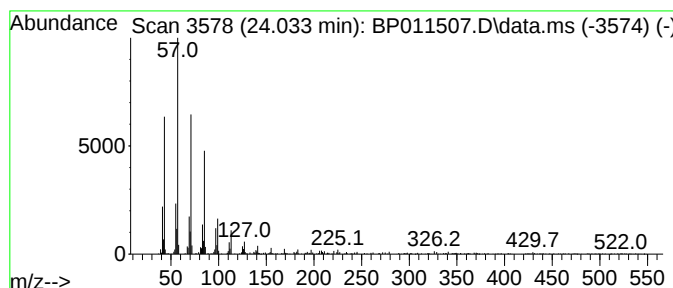
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.033	9.18 ng/ul	738023	Perylene-d12	23.433

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	95
2		Nonadecane	268	C19H40	000629-92-5	95
3		Tetracosane	338	C24H50	000646-31-1	93
4		Heptadecane	240	C17H36	000629-78-7	93
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011507.D
Acq On : 19 Aug 2022 14:53
Operator : CG/JU
Sample : N4210-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

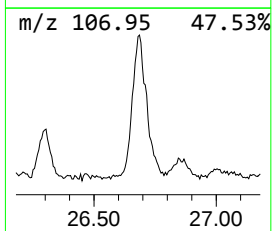
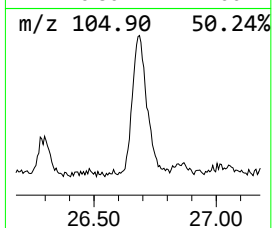
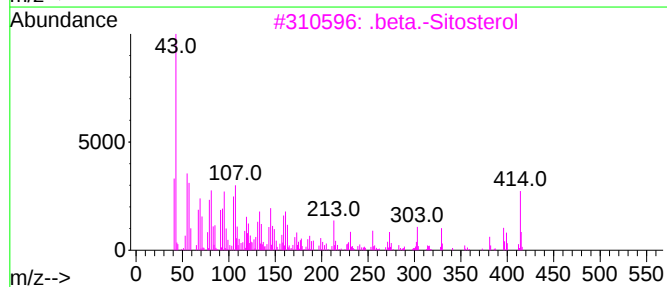
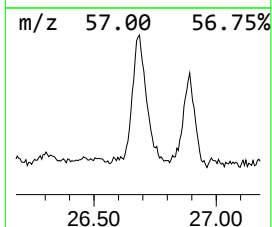
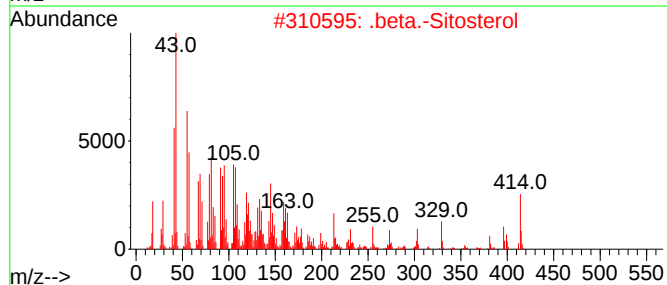
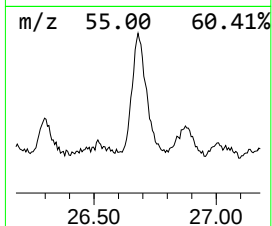
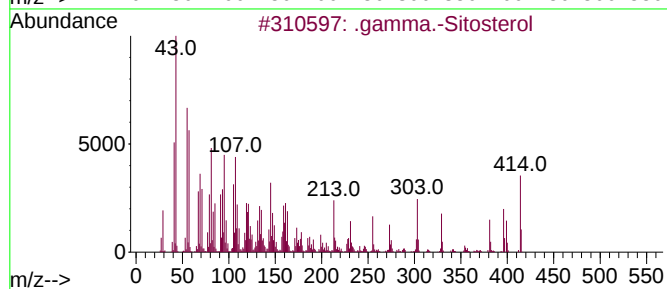
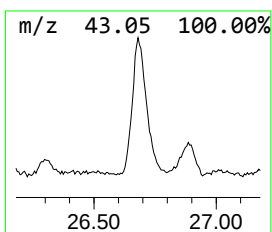
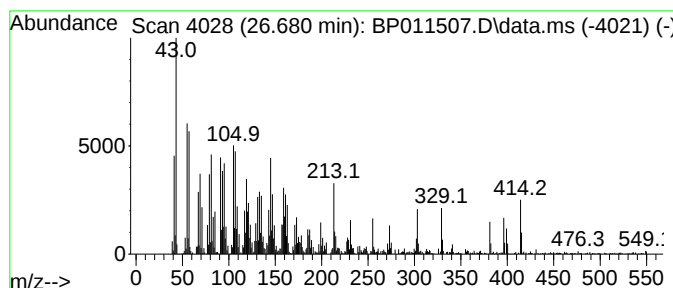
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 .gamma.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.680	29.99 ng/ul	2410270	Perylene-d12	23.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	98
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	95
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	94
4	.gamma.-Sitosterol	414	C29H50O	000083-47-6	83
5	(3S,8S,9S,10R,13R,14S,17R)-17-((...	414	C29H50O	029365-29-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081922\
Data File : BP011507.D
Acq On : 19 Aug 2022 14:53
Operator : CG/JU
Sample : N4210-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BGC49

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Palmitoleic acid	17.869	4.2	ng/ul	323768	4	17.004	1546870	20.0
n-Hexadecanoic ...	17.928	7.7	ng/ul	594689	4	17.004	1546870	20.0
(DEL) Alkane: S...	20.869	5.2	ng/ul	526896	5	21.133	2015210	20.0
(DEL) Alkane: S...	21.739	2.8	ng/ul	277079	5	21.133	2015210	20.0
Terephthalic ac...	22.004	4.3	ng/ul	431377	5	21.133	2015210	20.0
(DEL) Alkane: S...	24.033	9.2	ng/ul	738023	6	23.433	1607200	20.0
.gamma.-Sitosterol	26.680	30.0	ng/ul	2410270	6	23.433	1607200	20.0