

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
 Data File : BN020035.D
 Acq On : 07 Jun 2022 19:16
 Operator : CG/JU
 Sample : N3171-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBTZ6

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_N\Methods\SFAM-EPA-BN060622.M
 Title : SVOA CALIBRATION

Signal : TIC: BN020035.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	30	34	42	rVB	20787	33387	0.83%	0.086%
2	3.552	76	79	91	rBV	28521	49017	1.22%	0.126%
3	3.699	101	104	113	rBV	38014	66310	1.65%	0.170%
4	3.934	140	144	150	rVB3	9001	15660	0.39%	0.040%
5	4.028	156	160	164	rVB	8837	11598	0.29%	0.030%
6	4.246	192	197	203	rBV4	12331	20209	0.50%	0.052%
7	4.658	263	267	271	rBV2	7775	10220	0.25%	0.026%
8	4.940	309	315	321	rBV7	3110	6774	0.17%	0.017%
9	5.534	413	416	422	rVB7	2404	5056	0.13%	0.013%
10	5.922	477	482	488	rVB6	3696	6565	0.16%	0.017%
11	6.228	529	534	540	rBV2	2944	5628	0.14%	0.014%
12	6.817	628	634	638	rBV5	3156	6189	0.15%	0.016%
13	6.987	657	663	671	rBV	91294	155674	3.88%	0.399%
14	7.152	683	691	699	rBV	401033	659772	16.43%	1.693%
15	7.222	699	703	712	rVB	38258	65324	1.63%	0.168%
16	7.352	718	725	736	rBV	367182	618442	15.40%	1.587%
17	7.822	798	805	814	rBV	461532	763598	19.01%	1.959%
18	7.893	814	817	821	rVB2	8167	12146	0.30%	0.031%
19	8.528	916	925	935	rBV	235854	399605	9.95%	1.025%
20	8.916	986	991	994	rBV	6180	9298	0.23%	0.024%
21	8.975	994	1001	1015	rVV	508256	877067	21.84%	2.251%
22	9.099	1015	1022	1026	rVV5	7838	14451	0.36%	0.037%
23	9.152	1026	1031	1038	rVB6	5743	11129	0.28%	0.029%
24	9.434	1071	1079	1085	rVB2	5928	11329	0.28%	0.029%
25	9.693	1116	1123	1131	rBV	386861	669416	16.67%	1.718%
26	9.852	1144	1150	1154	rBV2	4370	7683	0.19%	0.020%
27	10.234	1207	1215	1227	rBV	564552	970088	24.16%	2.489%
28	10.399	1239	1243	1248	rBV5	3916	7466	0.19%	0.019%
29	10.610	1271	1279	1290	rBV	666957	1195516	29.77%	3.068%
30	10.746	1294	1302	1312	rBV	93822	177942	4.43%	0.457%
31	11.922	1499	1502	1510	rVB6	3492	5410	0.13%	0.014%
32	12.057	1521	1525	1528	rVV5	3283	4914	0.12%	0.013%
33	12.204	1544	1550	1557	rVB2	10947	18847	0.47%	0.048%
34	12.816	1648	1654	1658	rBV6	2808	5806	0.14%	0.015%
35	13.069	1693	1697	1703	rBV7	3216	5992	0.15%	0.015%
36	13.293	1729	1735	1742	rBV2	14875	21108	0.53%	0.054%

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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_N\Methods\SFAM-EPA-BN060622.M
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37	13.357	1742	1746	1754	rVB5	8175	14768	0.37%	0.038%
38	13.863	1825	1832	1843	rBV	1263801	1874607	46.68%	4.810%
39	13.993	1850	1854	1858	rVB5	4226	6308	0.16%	0.016%
40	14.146	1870	1880	1887	rBV	1344696	2101916	52.34%	5.394%
41	14.363	1912	1917	1922	rVB2	12653	18056	0.45%	0.046%
42	14.451	1925	1932	1939	rBV	1138479	1645696	40.98%	4.223%
43	14.598	1954	1957	1959	rBV4	4514	5960	0.15%	0.015%
44	14.640	1959	1964	1976	rVV	112547	191599	4.77%	0.492%
45	14.798	1987	1991	1998	rVB8	6308	12270	0.31%	0.031%
46	14.869	1998	2003	2007	rVB5	4792	6918	0.17%	0.018%
47	14.928	2007	2013	2016	rVB7	3184	5559	0.14%	0.014%
48	14.981	2018	2022	2023	rBV4	4399	4872	0.12%	0.013%
49	15.004	2023	2026	2030	rVB4	5052	7459	0.19%	0.019%
50	15.263	2066	2070	2075	rBV2	8078	12684	0.32%	0.033%
51	15.316	2075	2079	2086	rVB	20479	28675	0.71%	0.074%
52	15.440	2093	2100	2113	rBV	1872013	2782127	69.28%	7.139%
53	15.557	2114	2120	2129	rBV	695491	976469	24.31%	2.506%
54	15.681	2136	2141	2145	rBV	21548	36751	0.92%	0.094%
55	15.716	2145	2147	2151	rVB3	7743	10966	0.27%	0.028%
56	15.875	2171	2174	2180	rVB9	5959	12171	0.30%	0.031%
57	15.940	2180	2185	2187	rBV5	4331	7789	0.19%	0.020%
58	16.175	2220	2225	2228	rBV	39859	56254	1.40%	0.144%
59	16.381	2258	2260	2267	rVB8	6265	6763	0.17%	0.017%
60	16.516	2281	2283	2286	rVB4	5408	4387	0.11%	0.011%
61	16.692	2308	2313	2316	rBV7	7114	16648	0.41%	0.043%
62	16.863	2338	2342	2348	rVB9	11463	19707	0.49%	0.051%
63	16.969	2355	2360	2365	rBV2	41558	68775	1.71%	0.176%
64	17.016	2366	2368	2372	rVB	13737	18597	0.46%	0.048%
65	17.192	2391	2398	2404	rBV	1358316	2028535	50.51%	5.205%
66	17.251	2404	2408	2409	rVV2	73854	87648	2.18%	0.225%
67	17.292	2409	2415	2426	rVB	2113240	3164353	78.79%	8.120%
68	17.428	2435	2438	2443	rBV6	19663	29302	0.73%	0.075%
69	17.575	2458	2463	2468	rBV6	40629	71084	1.77%	0.182%
70	17.704	2481	2485	2489	rBV	39166	52716	1.31%	0.135%
71	17.875	2509	2514	2520	rVB	172356	209179	5.21%	0.537%
72	18.092	2547	2551	2559	rBV	180294	283855	7.07%	0.728%
73	18.398	2600	2603	2612	rVB3	37954	72234	1.80%	0.185%
74	18.722	2656	2658	2661	rBV3	11384	11994	0.30%	0.031%
75	18.763	2663	2665	2670	rVB5	13128	14886	0.37%	0.038%
76	18.845	2676	2679	2686	rBV3	39216	54462	1.36%	0.140%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
 Title : SVOA CALIBRATION

77	19.051	2708	2714	2718	rVB2	184750	253624	6.32%	0.651%
78	19.151	2727	2731	2734	rBV3	28022	36050	0.90%	0.093%
79	19.181	2734	2736	2739	rVV2	28586	29481	0.73%	0.076%
80	19.222	2739	2743	2746	rVB	25797	32180	0.80%	0.083%
81	19.375	2765	2769	2780	rBV	92561	154151	3.84%	0.396%
82	19.581	2798	2804	2812	rBV	2596158	3819949	95.12%	9.802%
83	20.145	2897	2900	2906	rVB	31194	35930	0.89%	0.092%
84	21.098	3058	3062	3074	rBV	335461	586120	14.59%	1.504%
85	21.298	3094	3096	3103	rVB	20506	28111	0.70%	0.072%
86	21.375	3104	3109	3120	rVV	1746812	2315203	57.65%	5.941%
87	21.545	3134	3138	3148	rVB	65318	89531	2.23%	0.230%
88	21.992	3210	3214	3225	rVB	86339	150771	3.75%	0.387%
89	22.475	3292	3296	3302	rVB	105467	158882	3.96%	0.408%
90	23.010	3382	3387	3395	rBV	136326	223383	5.56%	0.573%
91	23.557	3472	3480	3487	rBV2	1909938	4015983	100.00%	10.305%
92	23.616	3487	3490	3497	rVV	151421	271162	6.75%	0.696%
93	23.710	3498	3506	3516	rVB2	1139968	2358296	58.72%	6.051%
94	24.310	3602	3608	3617	rVB	149524	310796	7.74%	0.798%
95	25.104	3734	3743	3754	rBV4	189345	655560	16.32%	1.682%
96	26.057	3900	3905	3919	rVB2	89913	283476	7.06%	0.727%
97	26.333	3947	3952	3969	rVB2	55126	202349	5.04%	0.519%

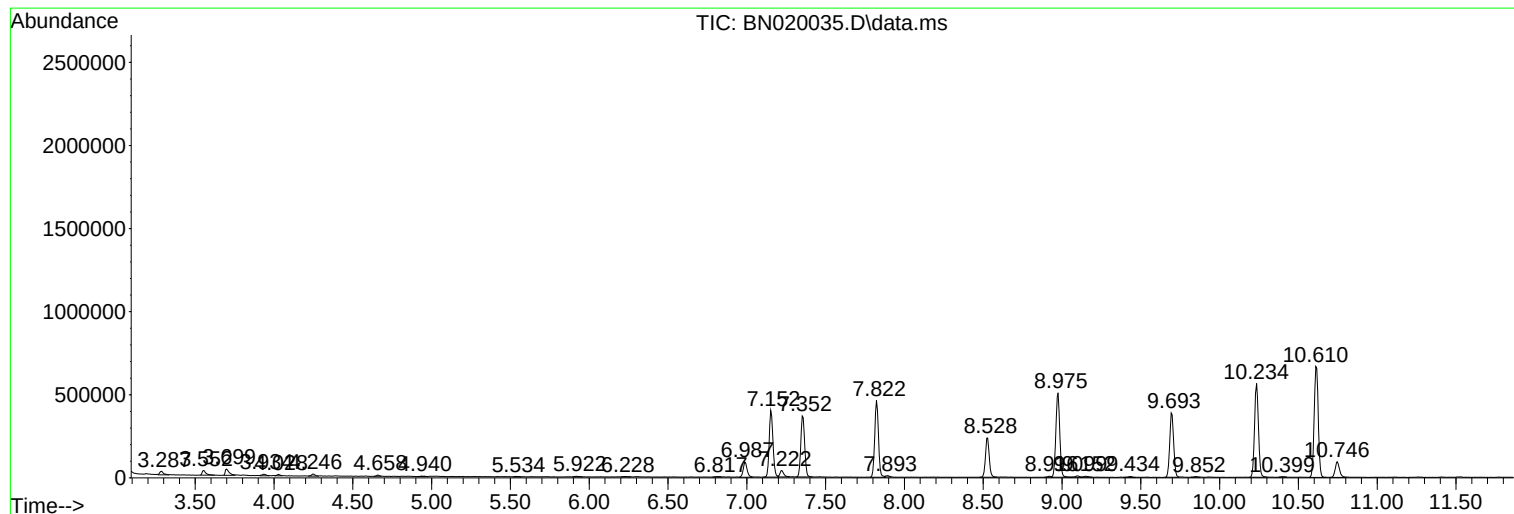
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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



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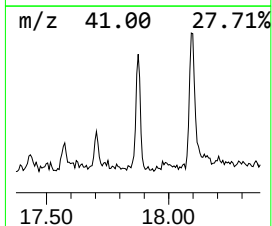
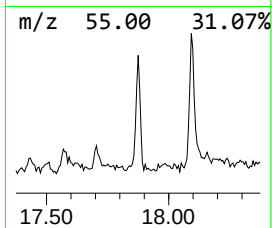
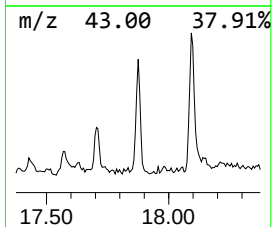
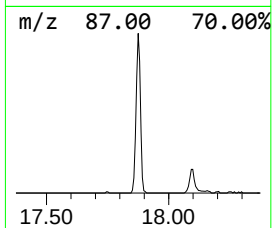
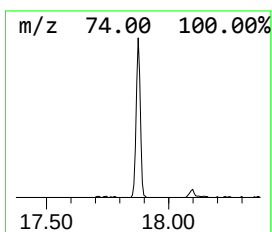
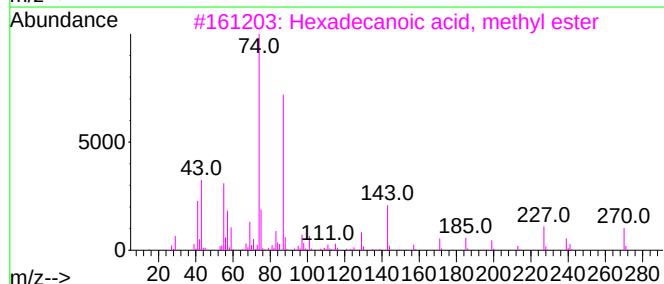
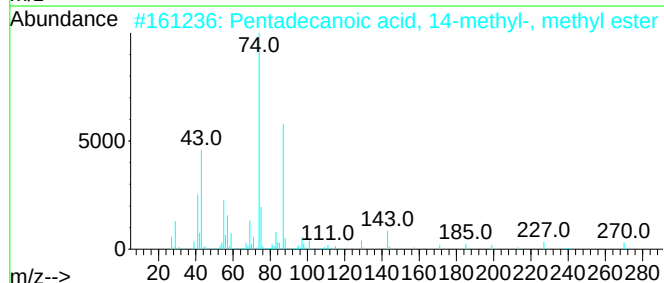
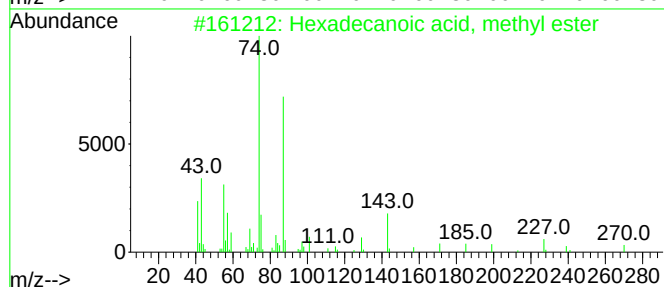
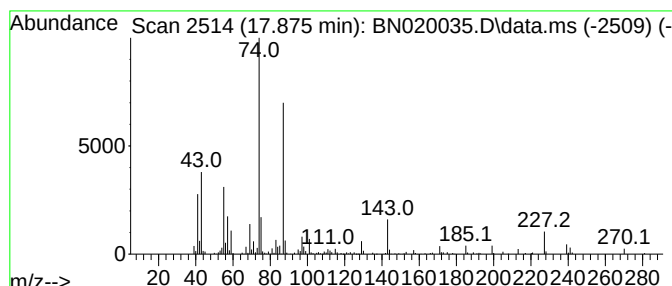
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Hexadecanoic acid, methyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.875	2.06 ng/ul	209179	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
5			Methyl 8-methyl-nonanoate	186	C11H22O2	1000452-00-9	89



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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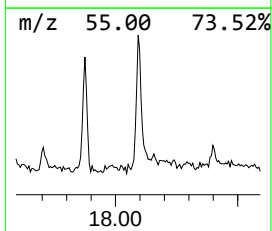
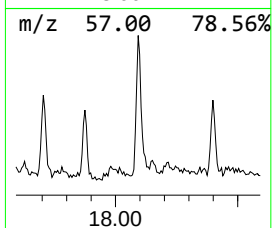
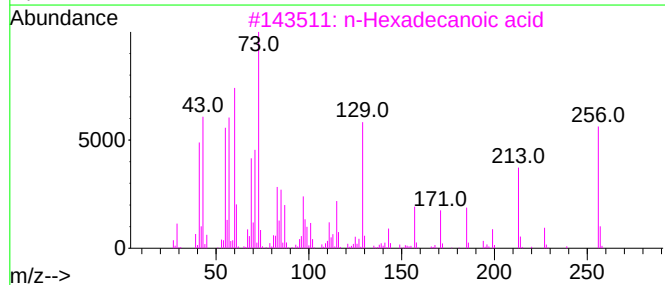
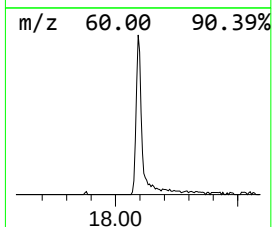
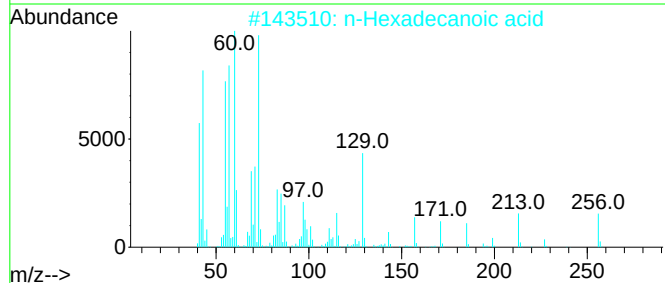
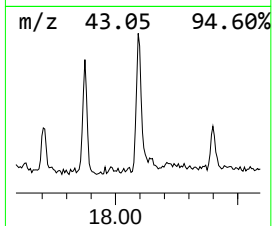
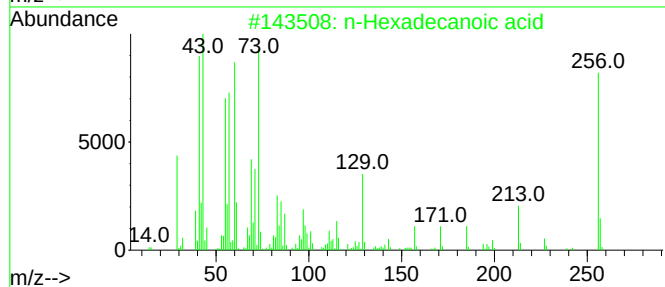
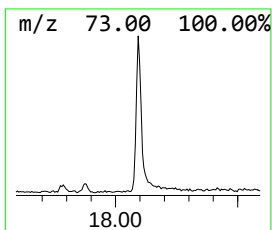
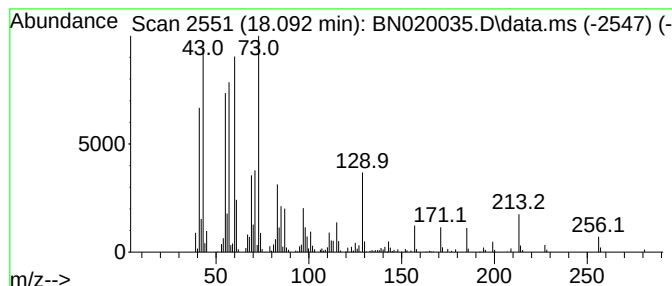
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	2.80 ng/ul	283855	Phenanthrene-d10	17.192

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



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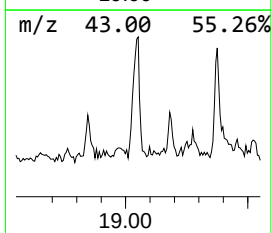
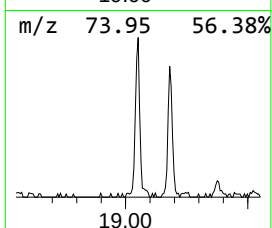
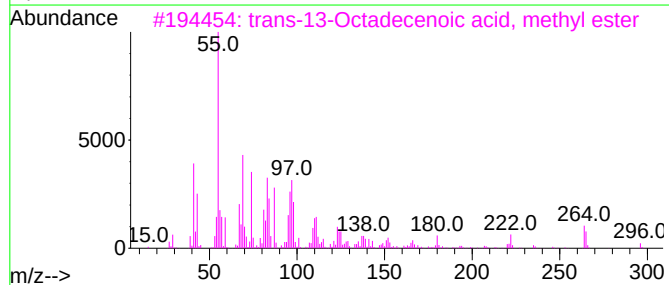
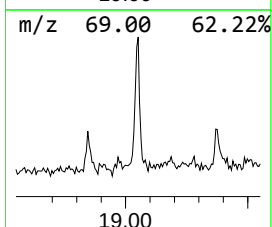
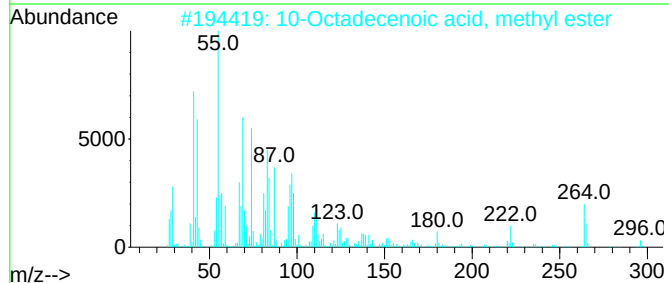
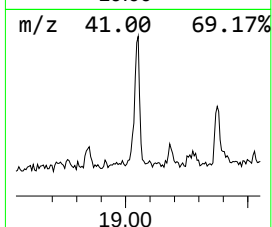
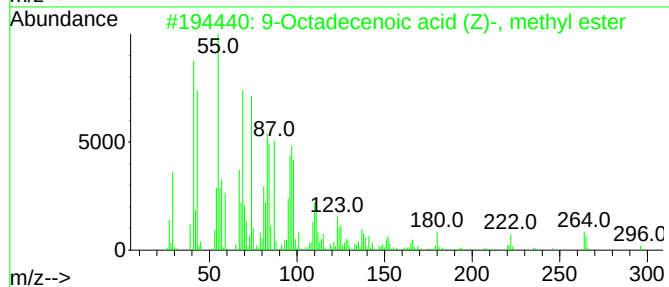
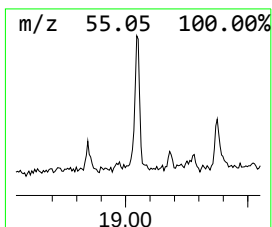
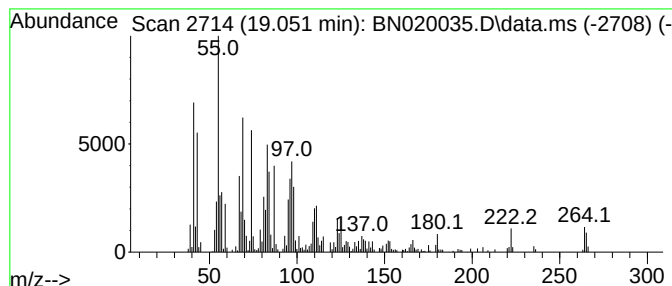
Instrument :
BNA_N
ClientSampleId :
DBTZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 9-Octadecenoic acid (Z)-, m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.051	2.50 ng/ul	253624	Phenanthrene-d10	17.192	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	95
2	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	94
3	trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	94
4	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	91
5	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	91



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ClientSampleId :
DBTZ6

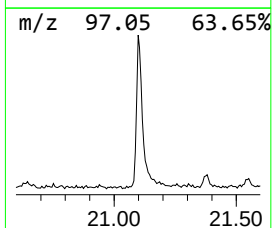
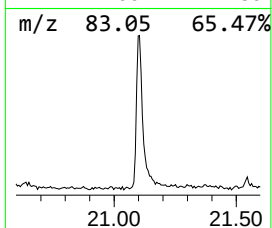
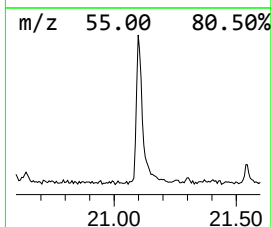
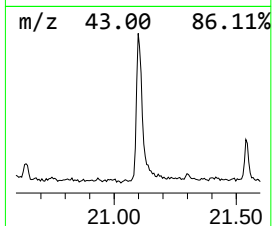
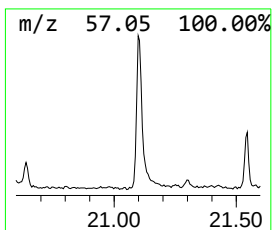
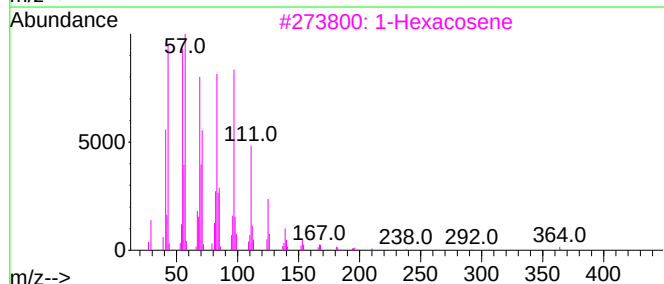
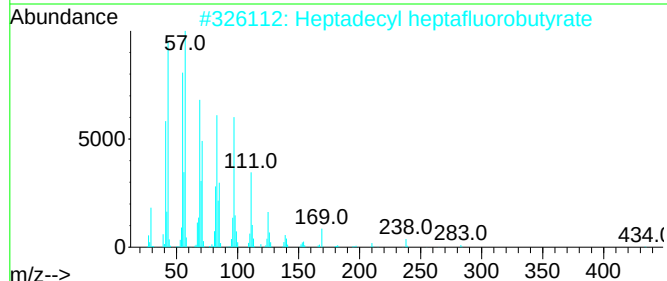
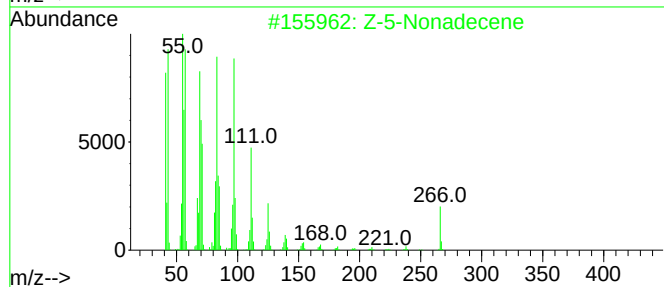
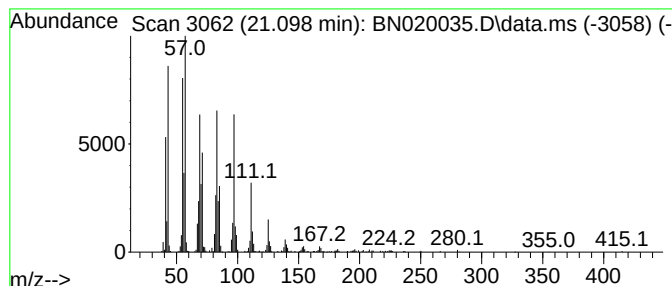
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	5.06 ng/ul	586120	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-5-Nonadecene	266	C19H38	1000131-11-8	96
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
5			Behenic alcohol	326	C22H46O	000661-19-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020035.D
Acq On : 07 Jun 2022 19:16
Operator : CG/JU
Sample : N3171-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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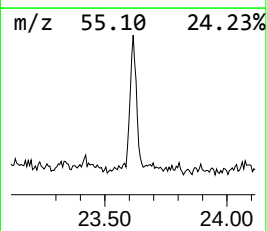
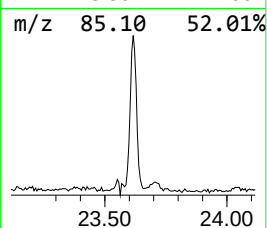
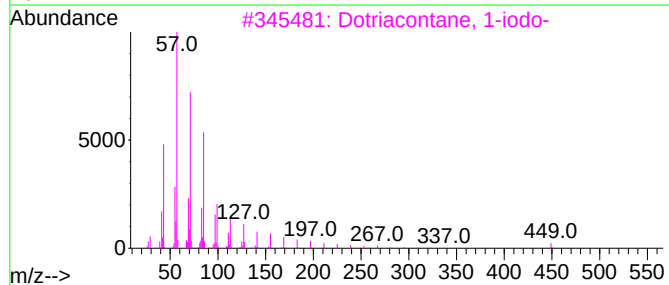
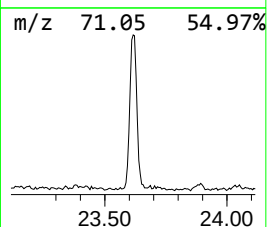
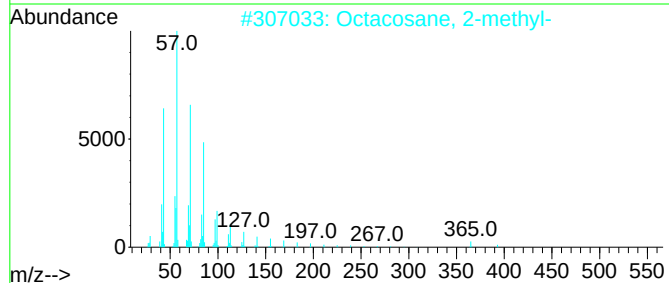
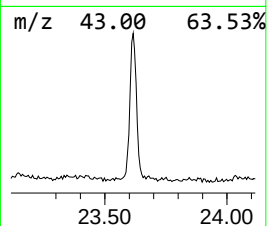
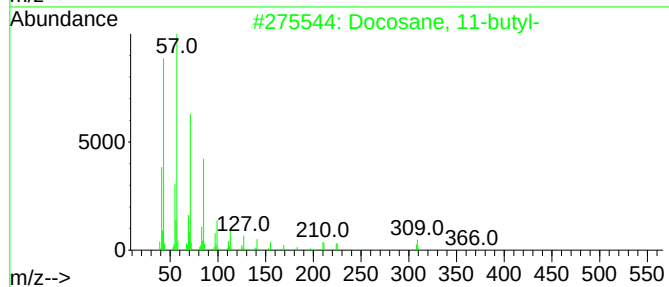
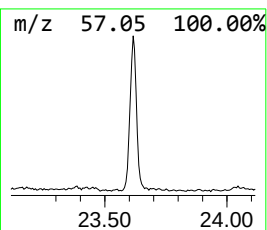
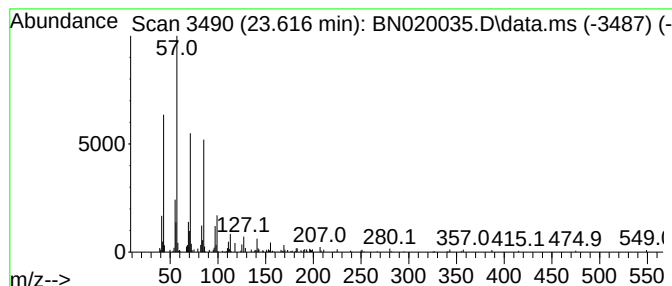
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.616	2.30 ng/ul	271162	Perylene-d12	23.710

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane, 11-butyl-	366	C26H54	013475-76-8	87
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	87
3		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	87
4		Tetracosane	338	C24H50	000646-31-1	87
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
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Operator : CG/JU
Sample : N3171-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

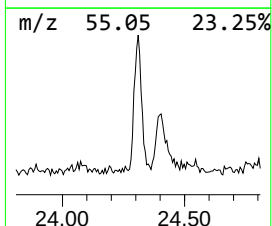
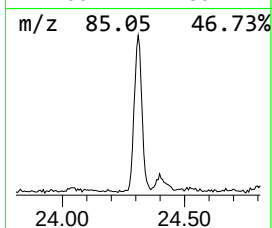
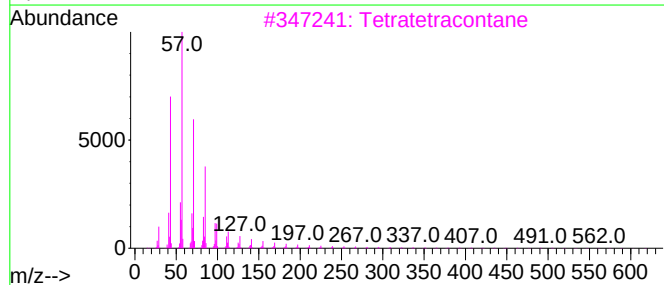
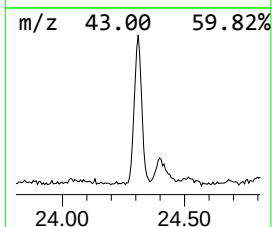
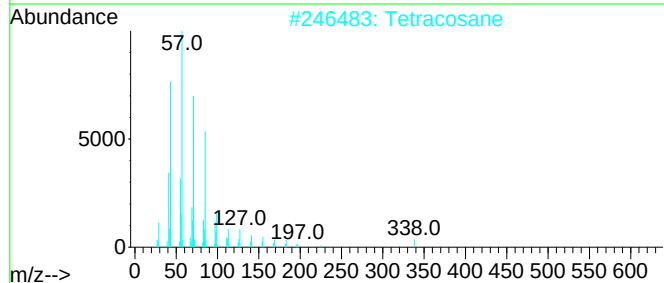
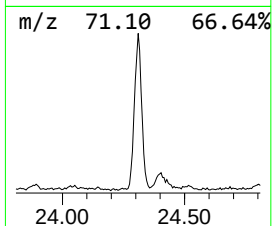
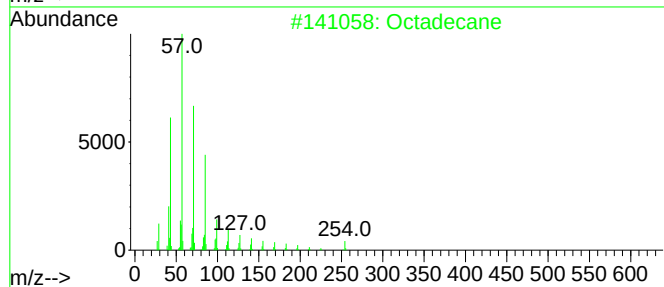
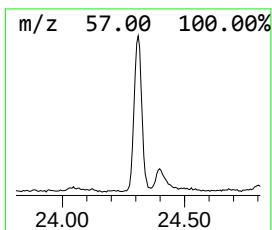
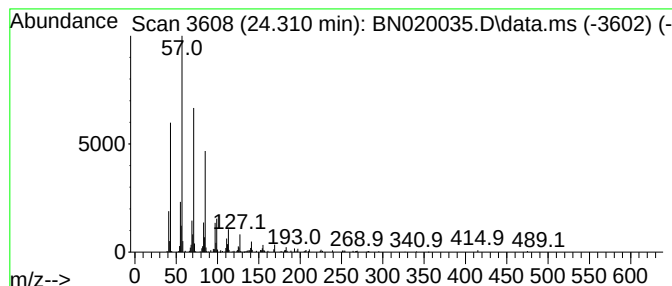
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.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.	
24.310	2.64 ng/ul	310796	Perylene-d12		23.710	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	94
2	Tetracosane		338	C24H50	000646-31-1	91
3	Tetratetracontane		619	C44H90	007098-22-8	91
4	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020035.D
Acq On : 07 Jun 2022 19:16
Operator : CG/JU
Sample : N3171-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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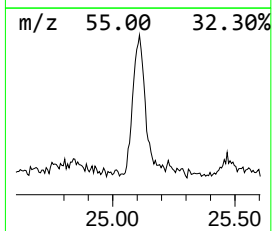
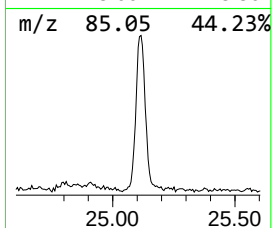
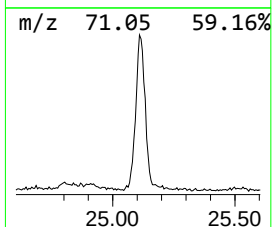
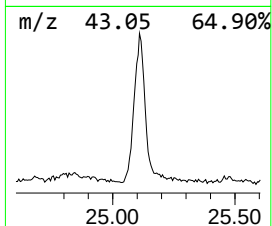
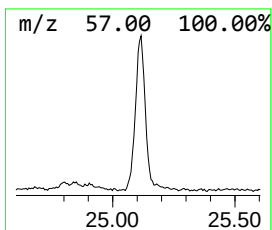
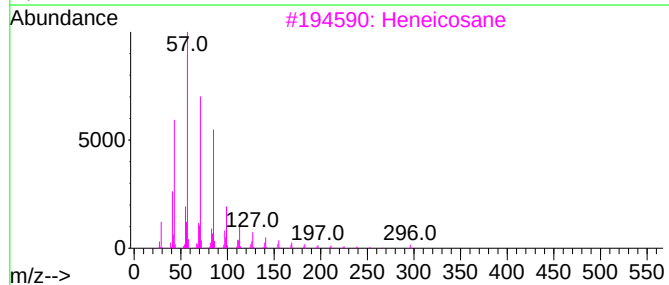
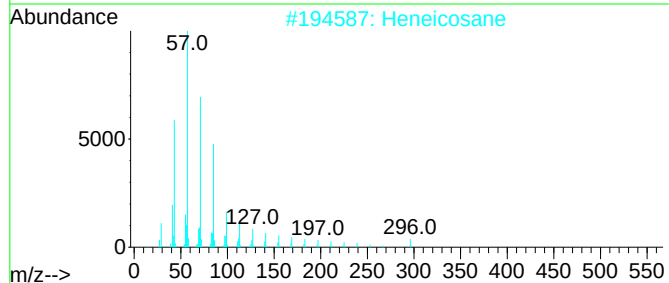
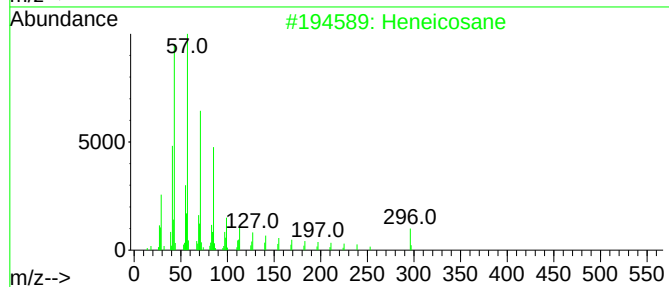
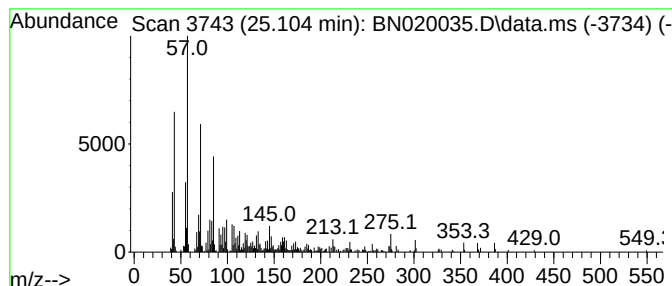
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.104	5.56 ng/ul	655560	Perylene-d12	23.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	92
2			Heneicosane	296	C21H44	000629-94-7	92
3			Heneicosane	296	C21H44	000629-94-7	91
4			26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	68
5			26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020035.D
Acq On : 07 Jun 2022 19:16
Operator : CG/JU
Sample : N3171-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :
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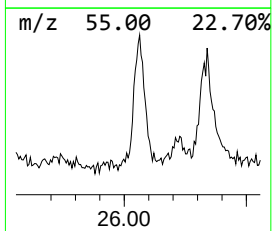
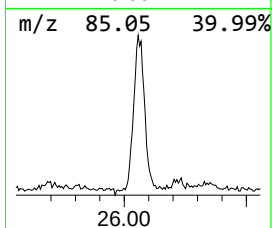
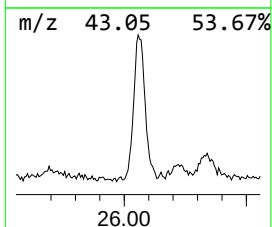
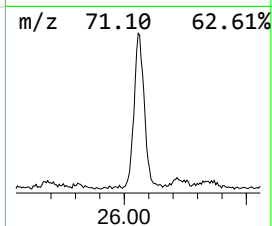
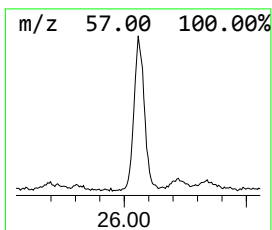
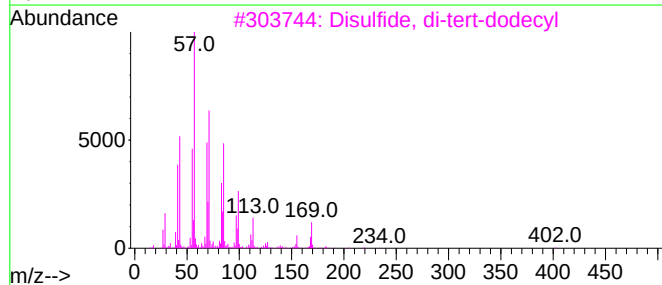
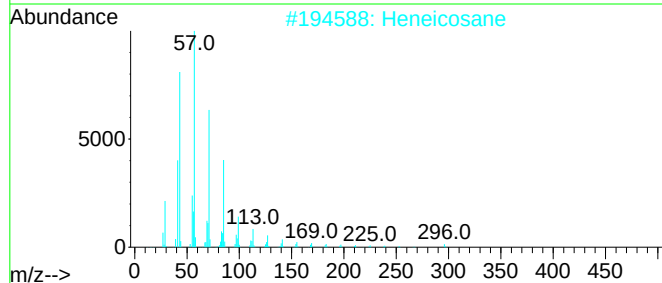
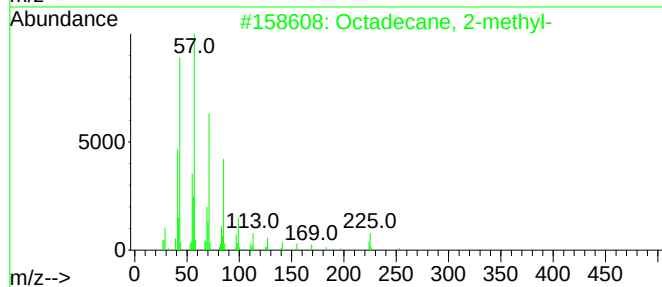
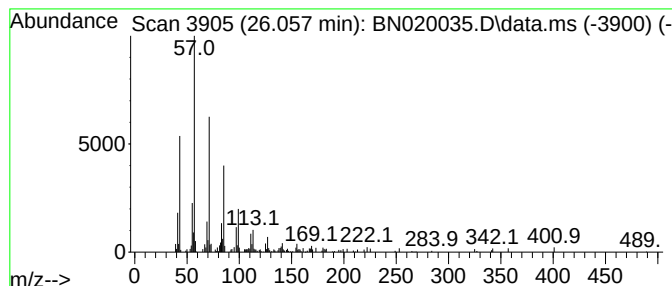
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.057	2.40 ng/ul	283476	Perylene-d12	23.710

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2-methyl-	268	C19H40	001560-88-9	86
2		Heneicosane	296	C21H44	000629-94-7	86
3		Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	86
4		Hentriacontane	437	C31H64	000630-04-6	83
5		11-Methylpentacosane	366	C26H54	015689-71-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060722\
Data File : BN020035.D
Acq On : 07 Jun 2022 19:16
Operator : CG/JU
Sample : N3171-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN060622.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic ...	18.092	2.8	ng/ul	283855	4	17.192	2028540	20.0
9-Octadecenoic ...	19.051	2.5	ng/ul	253624	4	17.192	2028540	20.0
(DEL) Alkane: S...	21.098	5.1	ng/ul	586120	5	21.375	2315200	20.0
(DEL) Alkane: S...	23.616	2.3	ng/ul	271162	6	23.710	2358300	20.0
(DEL) Alkane: S...	24.310	2.6	ng/ul	310796	6	23.710	2358300	20.0
(DEL) Alkane: S...	25.104	5.6	ng/ul	655560	6	23.710	2358300	20.0
(DEL) Alkane: S...	26.057	2.4	ng/ul	283476	6	23.710	2358300	20.0