

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
 Data File : BN023297.D
 Acq On : 20 Dec 2022 14:14
 Operator : CG/JU
 Sample : N6003-19
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBXT9

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-BN121522.M
 Title : SVOA CALIBRATION

Signal : TIC: BN023297.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	39	43	54	rVB	72527	114869	1.24%	0.101%
2	3.628	89	92	106	rVB	88444	134844	1.46%	0.118%
3	3.769	113	116	132	rBV	546168	914492	9.89%	0.803%
4	3.887	133	136	145	rVB	26107	39798	0.43%	0.035%
5	4.016	153	158	167	rVB	27741	48752	0.53%	0.043%
6	4.110	170	174	180	rVB	19443	26599	0.29%	0.023%
7	5.063	331	336	345	rVB2	131658	202631	2.19%	0.178%
8	6.969	655	660	667	rBV5	9569	21641	0.23%	0.019%
9	7.151	683	691	703	rVB	1403050	2190029	23.69%	1.923%
10	7.322	713	720	730	rVB	1139498	1728499	18.70%	1.518%
11	7.522	744	754	763	rBV	1459793	2235709	24.18%	1.963%
12	7.610	764	769	774	rVB5	8953	14363	0.16%	0.013%
13	7.998	827	835	842	rBV	1097230	1729018	18.70%	1.518%
14	8.057	842	845	853	rVB5	7564	13428	0.15%	0.012%
15	8.710	948	956	969	rBV	1577556	2500120	27.04%	2.195%
16	8.828	971	976	982	rVB	33936	53152	0.57%	0.047%
17	9.169	1025	1034	1042	rBV	1384104	2241658	24.25%	1.968%
18	9.328	1055	1061	1066	rVB9	7310	12381	0.13%	0.011%
19	9.581	1099	1104	1110	rVB2	23114	38592	0.42%	0.034%
20	9.892	1149	1157	1171	rBV	1206547	1960011	21.20%	1.721%
21	10.028	1176	1180	1185	rVB5	8178	13022	0.14%	0.011%
22	10.116	1190	1195	1201	rBV4	14140	21897	0.24%	0.019%
23	10.375	1234	1239	1242	rBV3	9195	14576	0.16%	0.013%
24	10.434	1242	1249	1259	rVB	1846269	3011219	32.57%	2.644%
25	10.592	1271	1276	1284	rVB2	28417	50025	0.54%	0.044%
26	10.816	1304	1314	1326	rBV	1667288	2756764	29.82%	2.421%
27	10.951	1326	1337	1354	rBV	1463401	2543284	27.51%	2.233%
28	11.604	1444	1448	1452	rBV5	7664	12328	0.13%	0.011%
29	12.075	1518	1528	1532	rVV6	6469	18663	0.20%	0.016%
30	12.228	1550	1554	1561	rVB	18134	25614	0.28%	0.022%
31	12.398	1578	1583	1589	rVB	35276	52642	0.57%	0.046%
32	13.433	1754	1759	1761	rBV3	13065	16606	0.18%	0.015%
33	13.469	1761	1765	1771	rVV	37620	52515	0.57%	0.046%
34	13.563	1777	1781	1785	rVV3	17146	30697	0.33%	0.027%
35	13.622	1785	1791	1797	rVB7	6497	17379	0.19%	0.015%
36	13.969	1845	1850	1854	rBV6	18006	25791	0.28%	0.023%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
 Title : SVOA CALIBRATION

37	14.057	1858	1865	1873	rVV	3311072	4429559	47.91%	3.889%
38	14.122	1873	1876	1880	rVB5	9725	13124	0.14%	0.012%
39	14.175	1880	1885	1890	rVB5	13466	23452	0.25%	0.021%
40	14.233	1890	1895	1900	rBV	14317	20161	0.22%	0.018%
41	14.339	1906	1913	1925	rVB	3886420	5551064	60.04%	4.874%
42	14.533	1943	1946	1951	rVB	22465	27724	0.30%	0.024%
43	14.645	1958	1965	1973	rBV	2730384	3874609	41.91%	3.402%
44	14.716	1973	1977	1983	rVB7	10850	18367	0.20%	0.016%
45	14.839	1991	1998	2012	rBV	1603146	2330783	25.21%	2.047%
46	14.992	2020	2024	2029	rVB3	22139	32515	0.35%	0.029%
47	15.057	2030	2035	2041	rVB3	49012	73921	0.80%	0.065%
48	15.639	2124	2134	2141	rBV	4783194	6816062	73.73%	5.985%
49	15.751	2147	2153	2160	rBV	1658553	2144534	23.20%	1.883%
50	15.874	2170	2174	2180	rVB2	26462	41476	0.45%	0.036%
51	16.004	2190	2196	2203	rVV	1507358	1990369	21.53%	1.748%
52	16.269	2237	2241	2246	rVB7	9887	13318	0.14%	0.012%
53	16.563	2286	2291	2297	rVV	32401	42716	0.46%	0.038%
54	16.786	2325	2329	2336	rBV6	16387	30547	0.33%	0.027%
55	17.092	2374	2381	2386	rBV	45846	70240	0.76%	0.062%
56	17.145	2386	2390	2393	rBV3	28377	35106	0.38%	0.031%
57	17.280	2410	2413	2416	rBV5	14837	21790	0.24%	0.019%
58	17.351	2421	2425	2426	rBV3	11968	14005	0.15%	0.012%
59	17.392	2426	2432	2438	rVV	3497498	4744508	51.32%	4.166%
60	17.492	2442	2449	2458	rBV2	5094692	7124565	77.06%	6.256%
61	17.580	2458	2464	2472	rVB6	45883	82438	0.89%	0.072%
62	17.645	2472	2475	2477	rBV2	13195	15977	0.17%	0.014%
63	17.769	2493	2496	2505	rVB3	30863	49923	0.54%	0.044%
64	18.016	2535	2538	2542	rBV3	13888	19773	0.21%	0.017%
65	18.057	2542	2545	2551	rVB3	26986	35044	0.38%	0.031%
66	18.110	2551	2554	2557	rBV4	11674	13437	0.15%	0.012%
67	18.168	2557	2564	2569	rBV4	33654	65567	0.71%	0.058%
68	18.227	2569	2574	2579	rBV	65382	103925	1.12%	0.091%
69	18.286	2579	2584	2594	rBV	1159455	1789063	19.35%	1.571%
70	18.533	2623	2626	2630	rBV4	21834	27520	0.30%	0.024%
71	18.586	2632	2635	2636	rVV3	16852	20043	0.22%	0.018%
72	18.616	2638	2640	2651	rVB3	40115	63829	0.69%	0.056%
73	18.710	2651	2656	2662	rBV	443767	504271	5.45%	0.443%
74	18.774	2663	2667	2676	rVB5	34573	60165	0.65%	0.053%
75	18.863	2680	2682	2689	rVB6	17707	20443	0.22%	0.018%
76	18.951	2695	2697	2703	rBV7	14542	24345	0.26%	0.021%

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77	19.139	2726	2729	2733	rBV4	32513	49312	0.53%	0.043%
78	19.186	2733	2737	2739	rBV5	33956	50464	0.55%	0.044%
79	19.351	2761	2765	2771	rBV2	82335	121569	1.31%	0.107%
80	19.415	2771	2776	2778	rBV2	186288	251416	2.72%	0.221%
81	19.445	2778	2781	2795	rVB2	186158	392871	4.25%	0.345%
82	19.563	2796	2801	2813	rBV	535506	830585	8.98%	0.729%
83	19.786	2833	2839	2852	rVB	6818038	9245163	100.00%	8.118%
84	19.945	2861	2866	2870	rBV	301271	377830	4.09%	0.332%
85	20.039	2878	2882	2885	rBV	175412	210415	2.28%	0.185%
86	20.298	2922	2926	2928	rBV3	55526	89748	0.97%	0.079%
87	20.657	2984	2987	2991	rVB2	118199	129778	1.40%	0.114%
88	20.698	2991	2994	3000	rBV4	79683	106583	1.15%	0.094%
89	20.815	3011	3014	3018	rVB3	107806	126391	1.37%	0.111%
90	21.274	3075	3092	3106	rVB2	2077425	5864309	63.43%	5.149%
91	21.580	3138	3144	3148	rBV	4015102	5999719	64.90%	5.268%
92	21.727	3166	3169	3178	rVB5	128620	212181	2.30%	0.186%
93	22.198	3245	3249	3250	rBV	244305	315196	3.41%	0.277%
94	23.262	3426	3430	3436	rBV2	487028	771608	8.35%	0.678%
95	23.868	3526	3533	3548	rVB2	3976409	8672530	93.81%	7.615%
96	24.027	3553	3560	3570	rVB2	2301176	4738529	51.25%	4.161%
97	24.651	3660	3666	3675	rVB	829550	1736605	18.78%	1.525%
98	24.750	3678	3683	3697	rVB4	293228	791542	8.56%	0.695%
99	26.539	3979	3987	4004	rVB	818774	2624241	28.39%	2.304%
100	29.086	4411	4420	4433	rVB6	800003	2913734	31.52%	2.558%

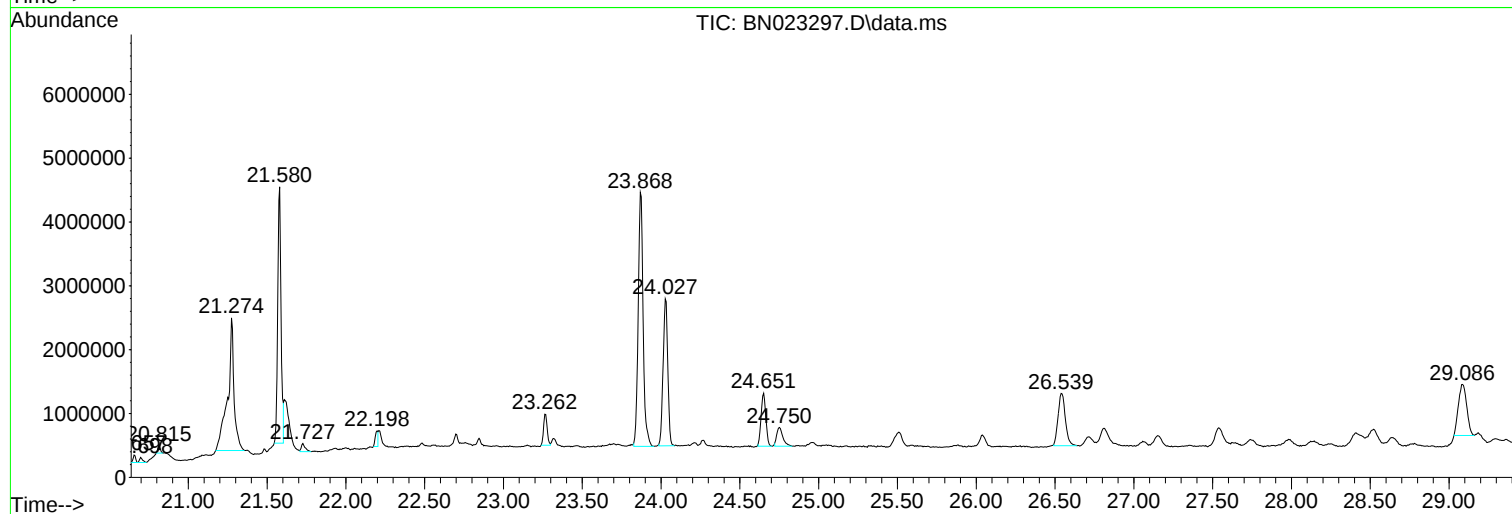
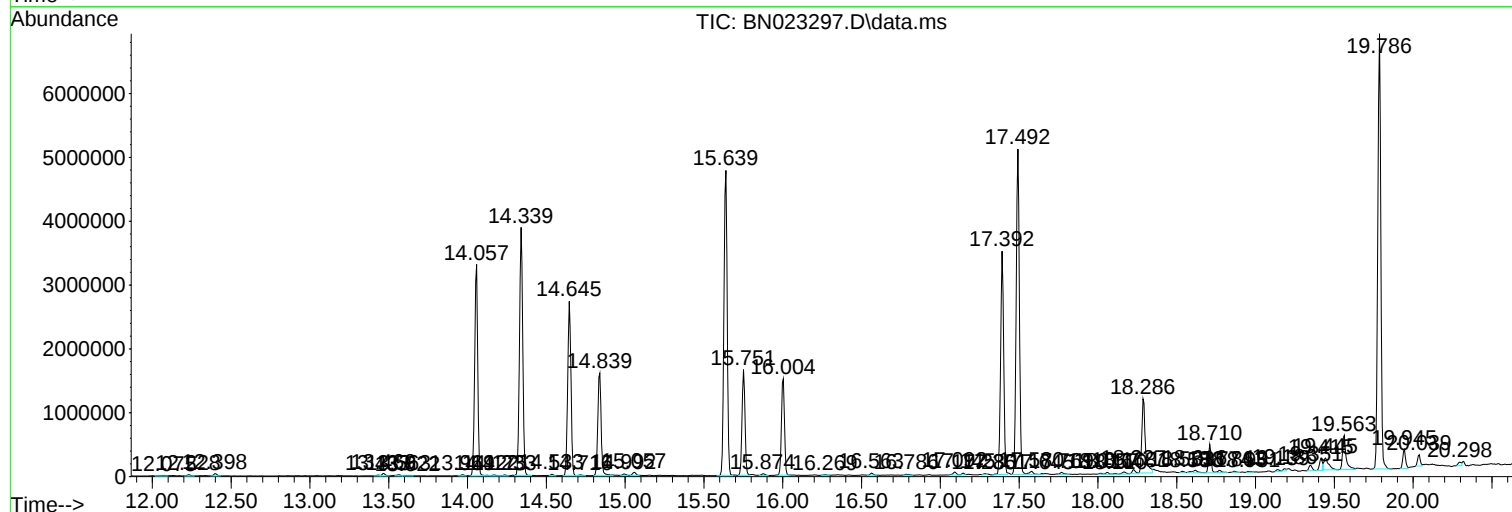
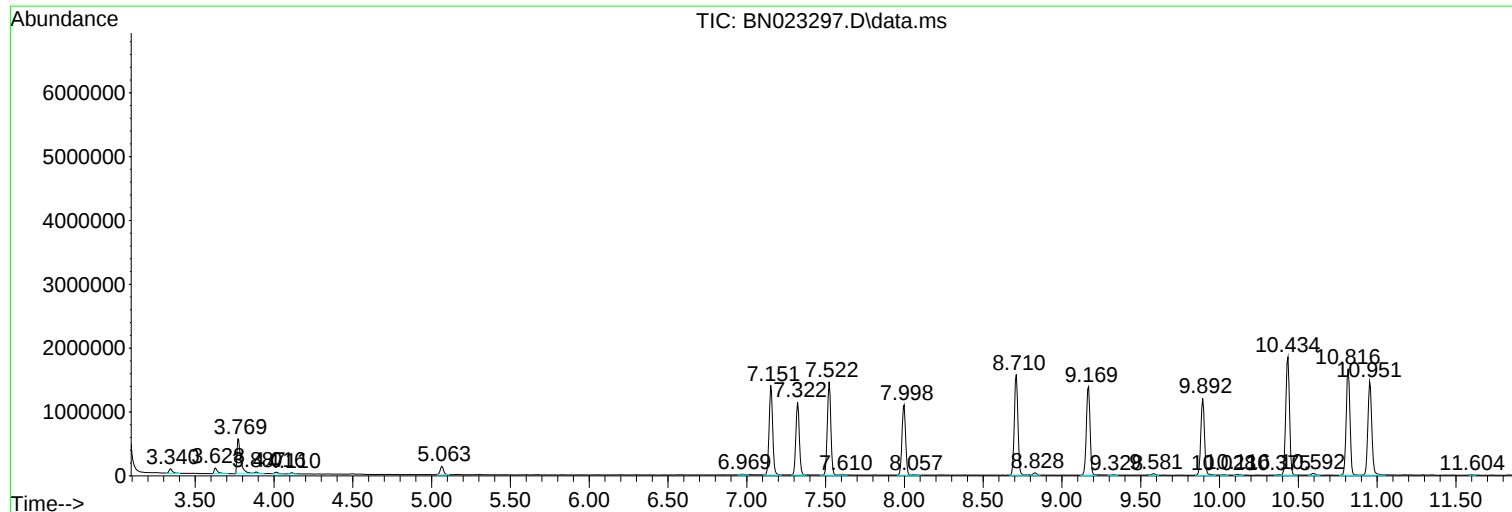
Sum of corrected areas: 113886205

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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



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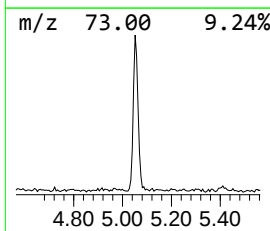
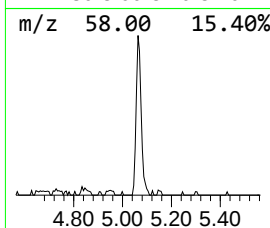
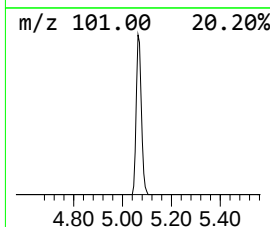
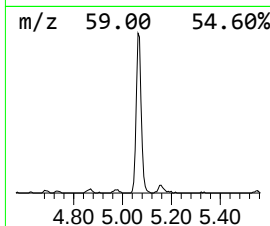
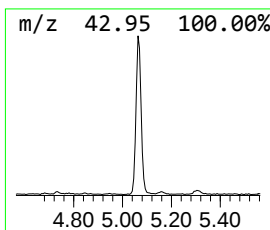
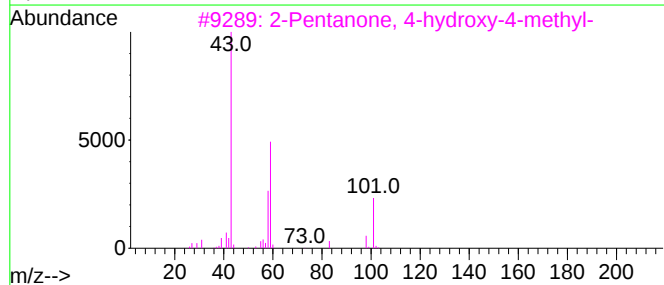
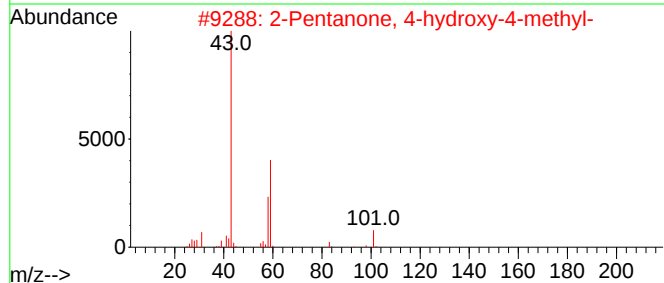
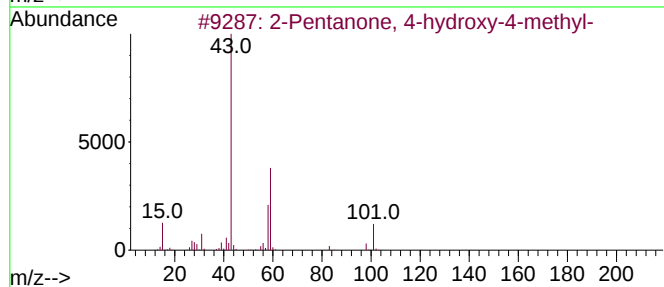
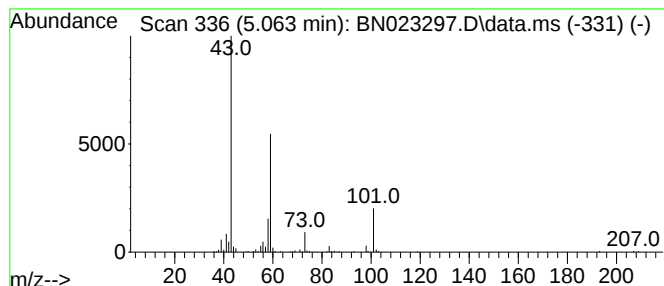
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.063	2.34 ng/ul	202631	1,4-Dichlorobenzene-d4	7.998

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25		



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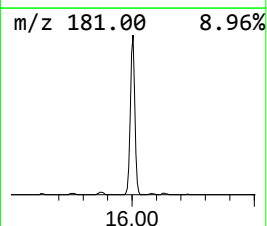
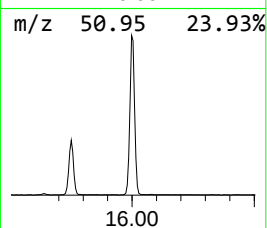
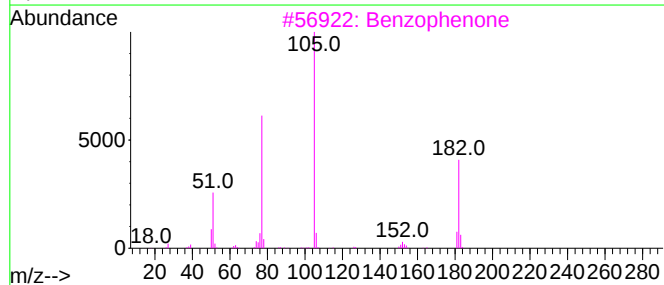
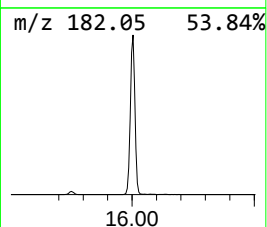
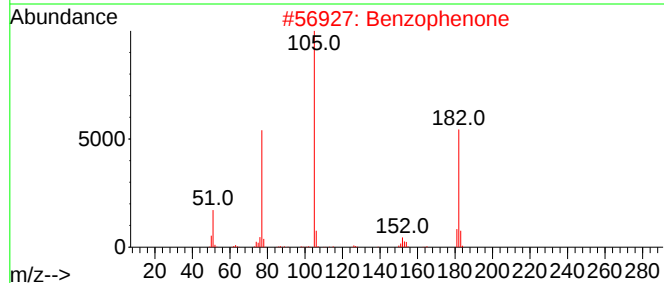
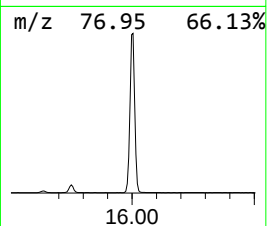
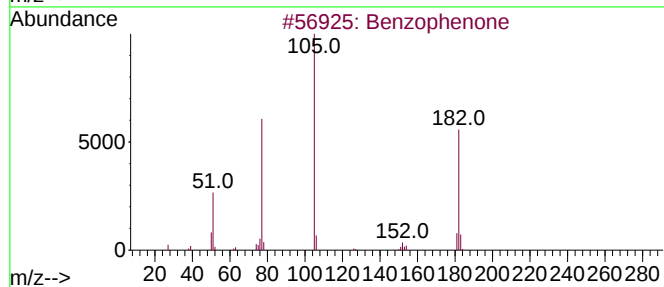
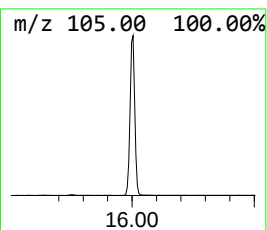
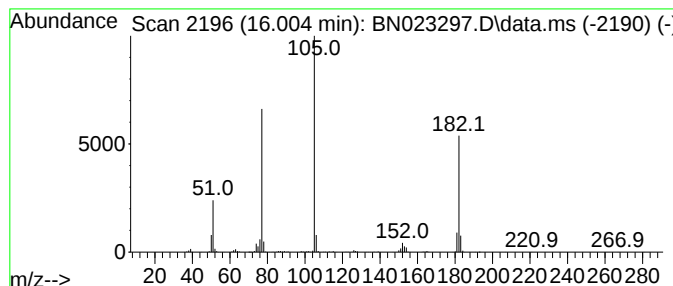
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.004	10.27 ng/ul	1990370	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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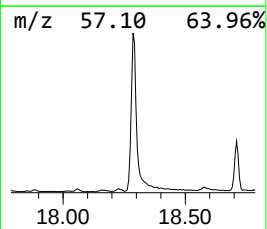
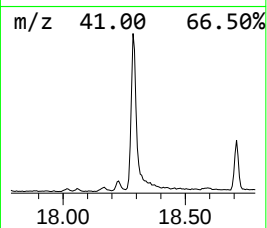
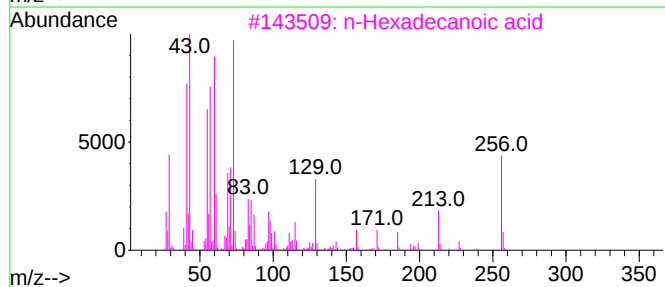
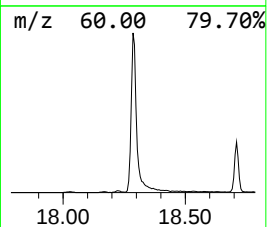
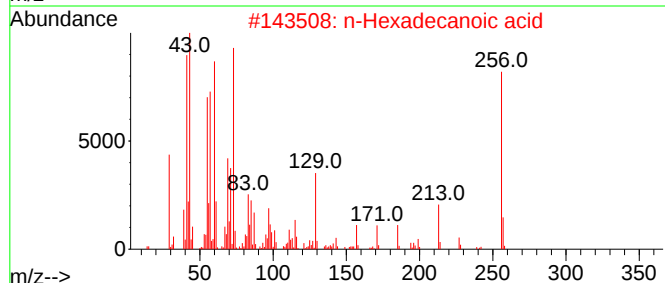
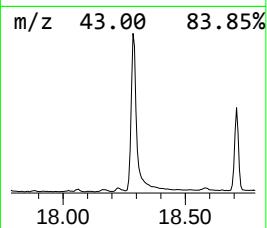
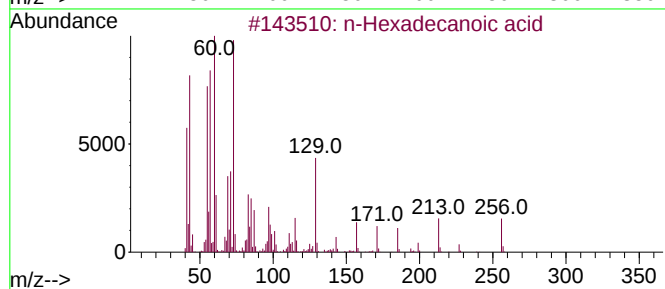
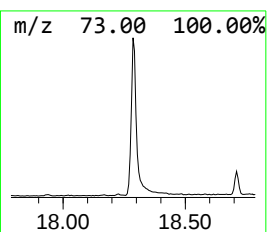
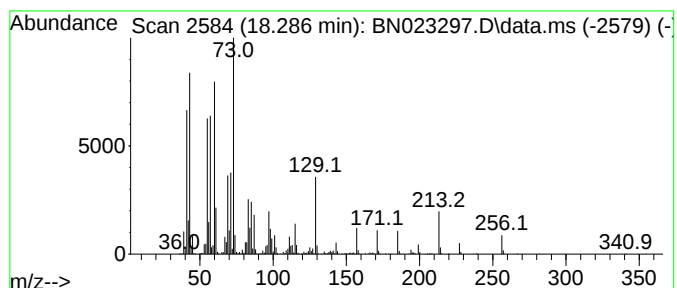
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ClientSampleId :
DBXT9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.286	7.54 ng/ul	1789060	Phenanthrene-d10	17.392	
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	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
Operator : CG/JU
Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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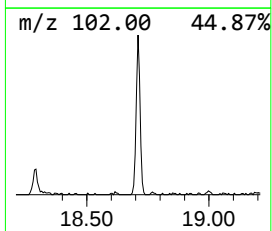
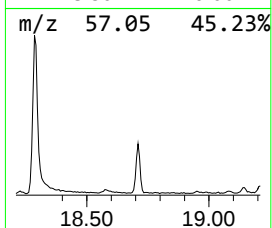
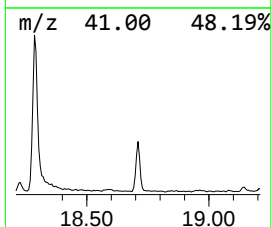
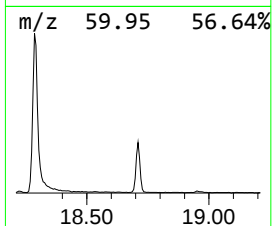
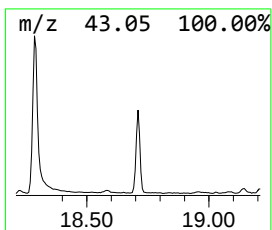
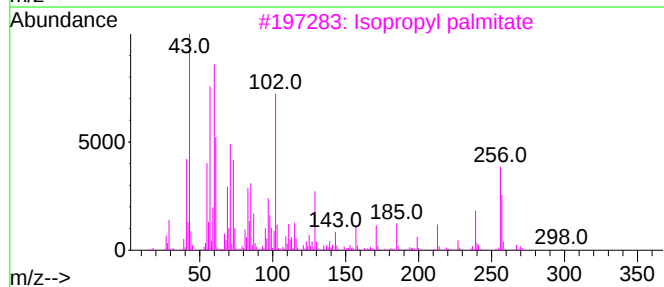
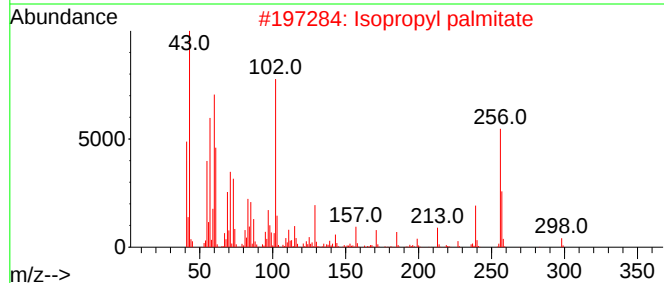
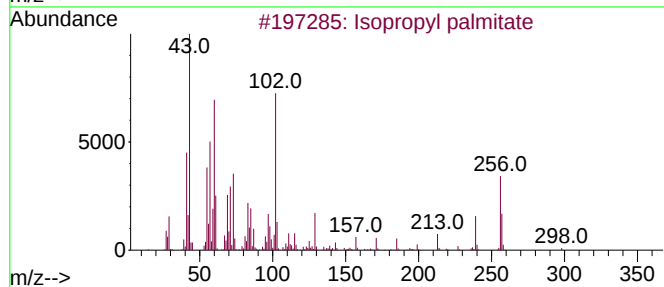
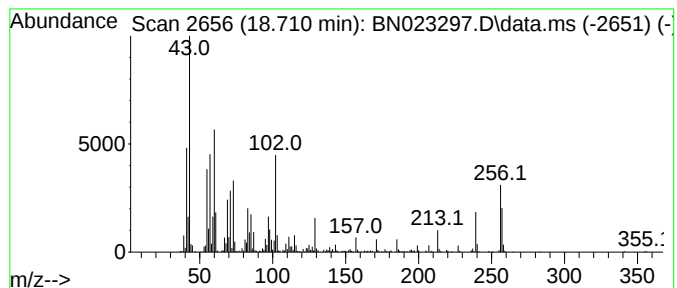
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Isopropyl palmitate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.710	2.13 ng/ul	504271	Phenanthrene-d10	17.392

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl palmitate	298	C19H38O2	000142-91-6	91
2			Isopropyl palmitate	298	C19H38O2	000142-91-6	86
3			Isopropyl palmitate	298	C19H38O2	000142-91-6	74
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	62
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
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Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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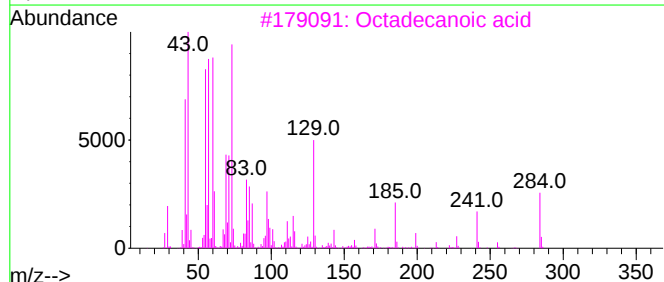
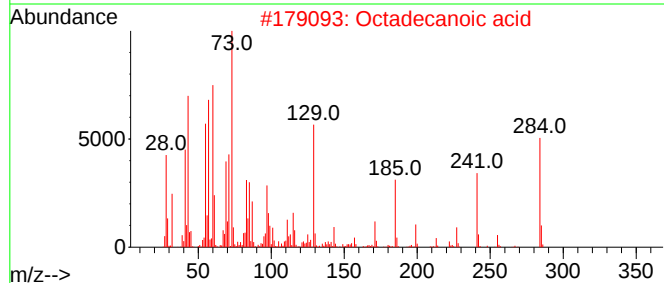
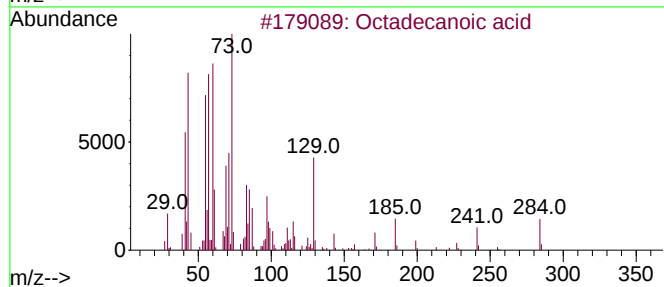
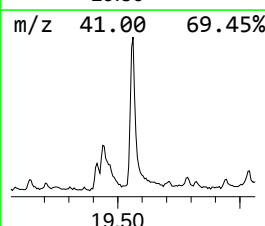
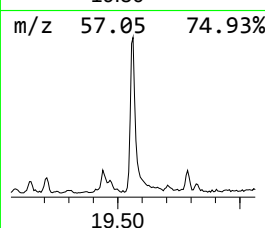
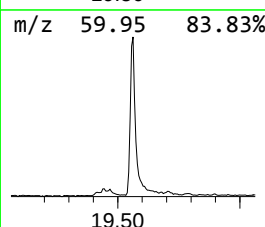
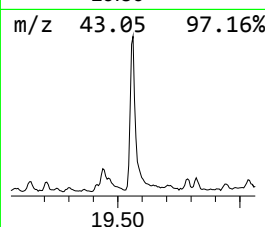
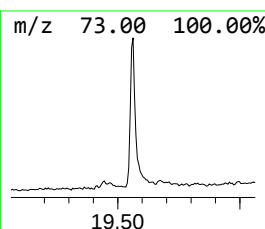
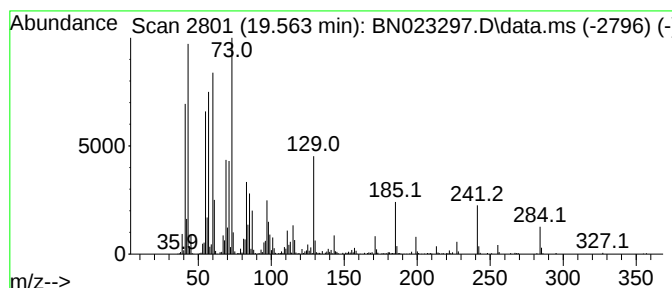
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.563	2.77 ng/ul	830585	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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	99
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
Operator : CG/JU
Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

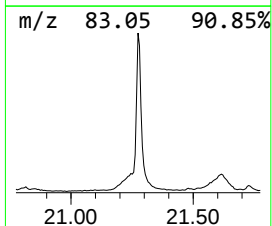
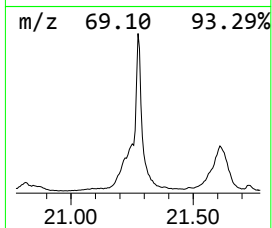
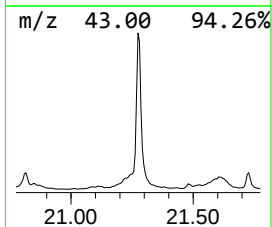
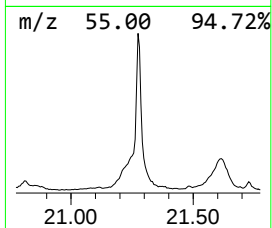
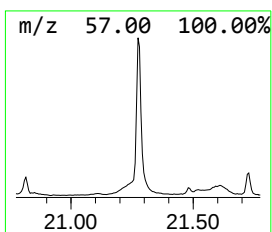
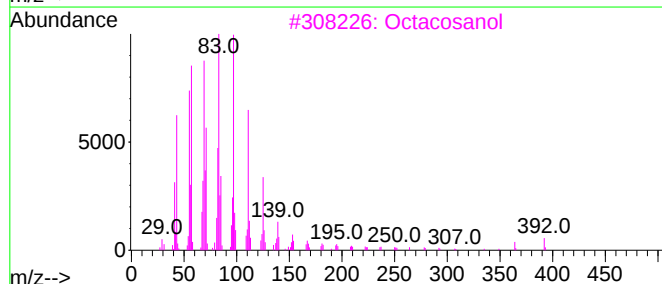
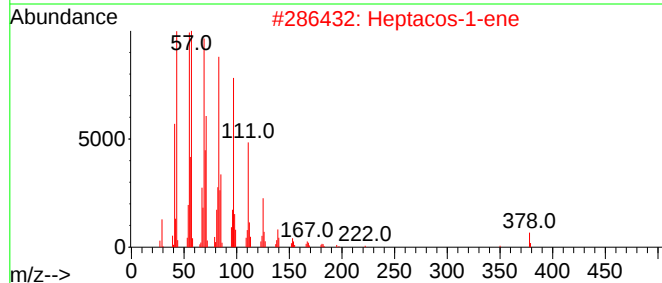
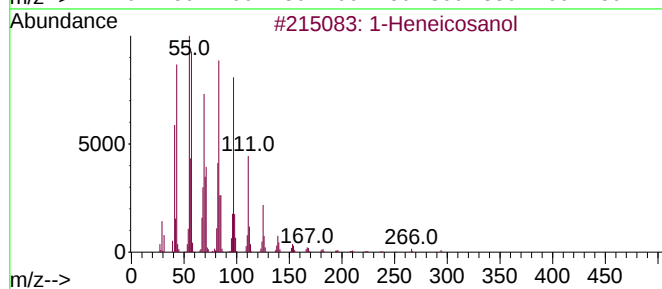
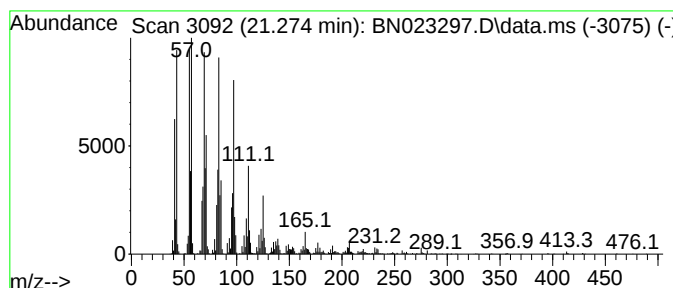
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.274	19.55 ng/ul	5864310	Chrysene-d12	21.580	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	93
2	Heptacos-1-ene	378	C27H54	015306-27-1	93
3	Octacosanol	410	C28H58O	000557-61-9	91
4	Cyclopentadecane	210	C15H30	000295-48-7	91
5	1-Heptacosanol	396	C27H56O	002004-39-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
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Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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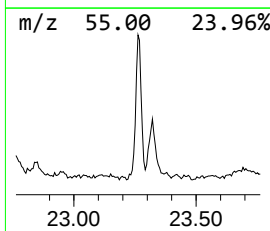
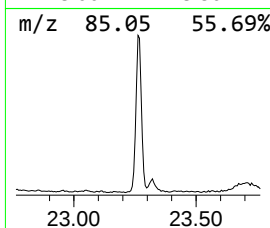
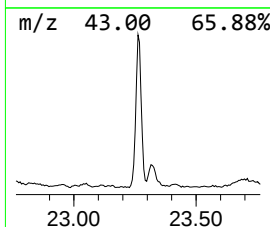
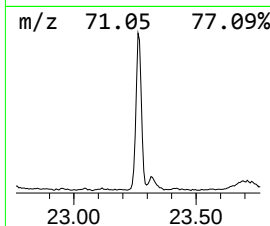
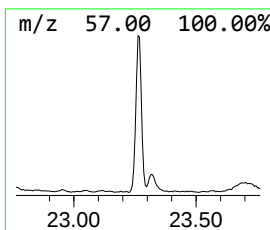
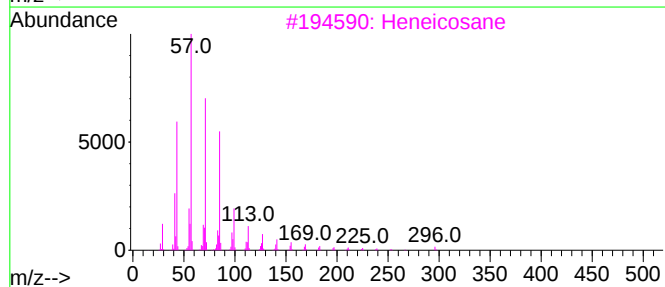
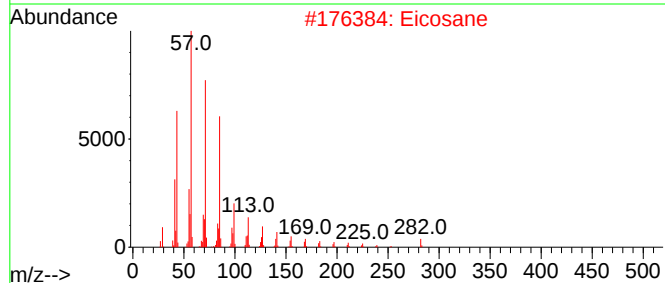
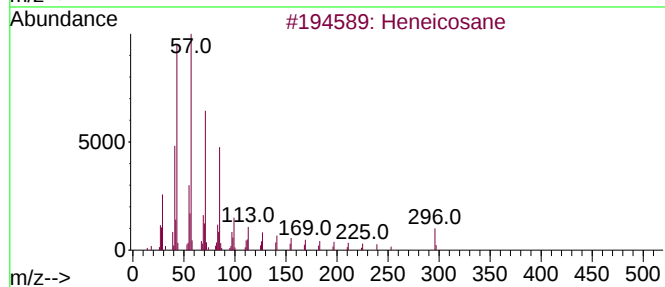
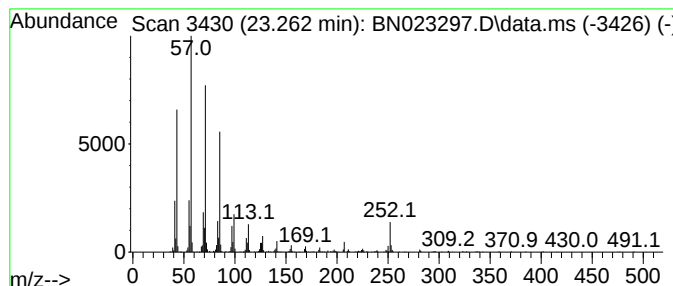
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.262	3.26 ng/ul	771608	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	93
2			Eicosane	282	C20H42	000112-95-8	93
3			Heneicosane	296	C21H44	000629-94-7	90
4			Hexacosane	366	C26H54	000630-01-3	90
5			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
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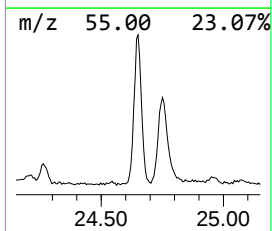
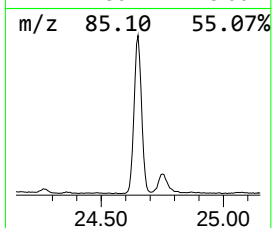
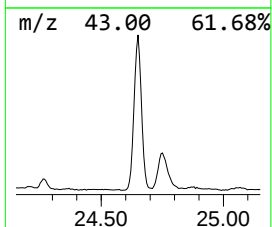
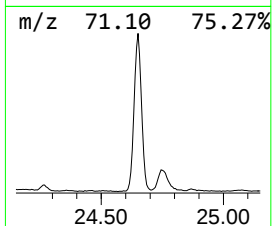
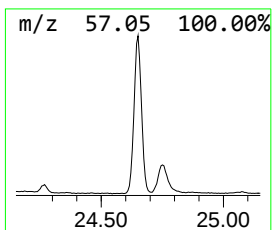
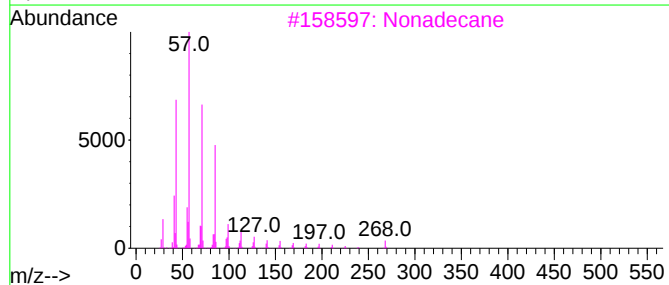
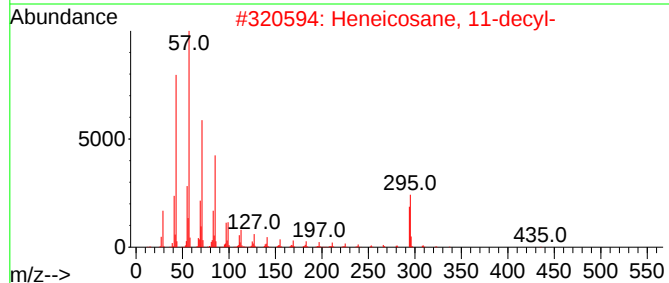
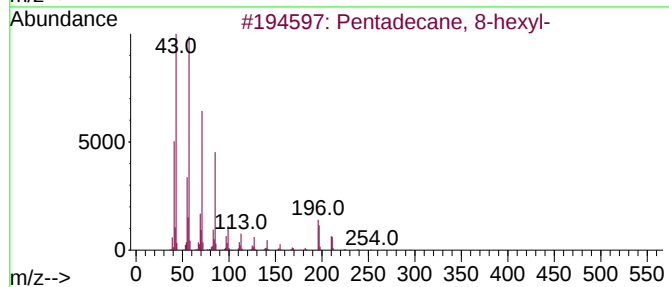
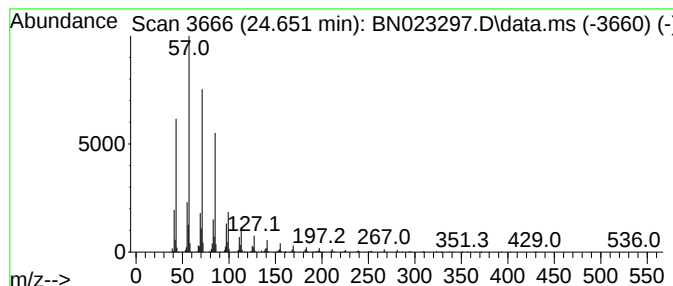
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Quant Title : SVOA CALIBRATION

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.
24.651	7.33 ng/ul	1736610	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	94
3		Nonadecane	268	C19H40	000629-92-5	93
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
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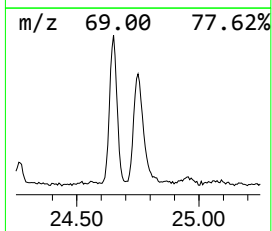
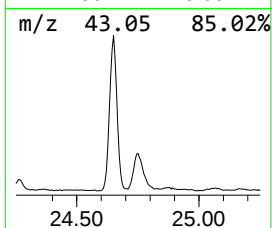
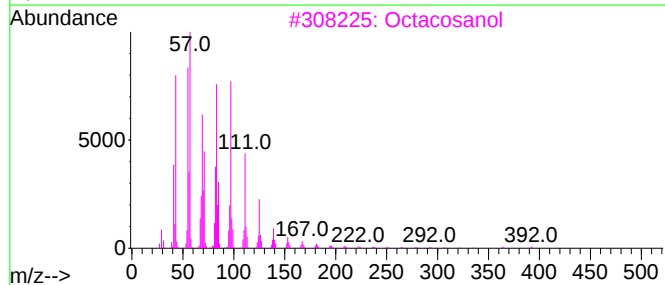
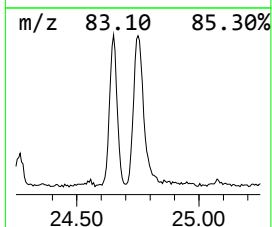
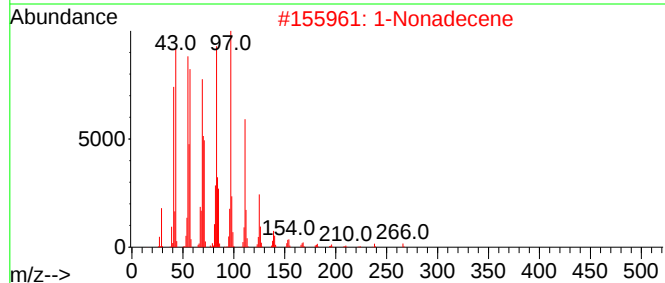
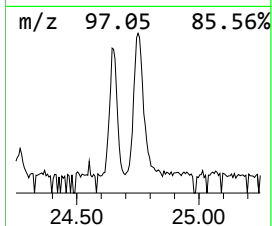
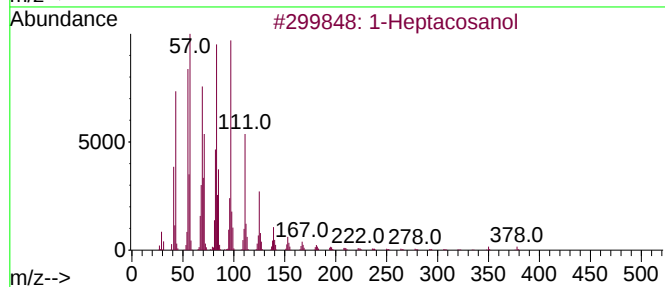
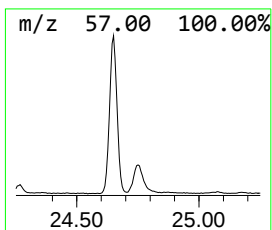
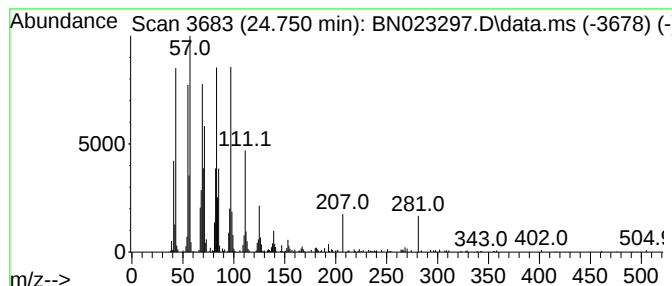
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heptacosanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.750	3.34 ng/ul	791542	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptacosanol	396	C27H56O	002004-39-9	95
2			1-Nonadecene	266	C19H38	018435-45-5	94
3			Octacosanol	410	C28H58O	000557-61-9	93
4			Nonacos-1-ene	406	C29H58	018835-35-3	93
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
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Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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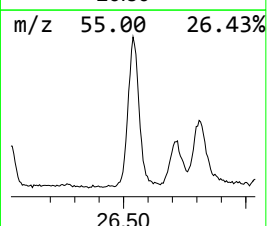
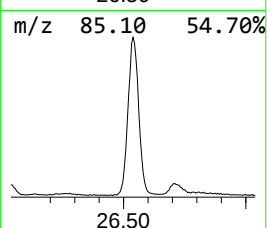
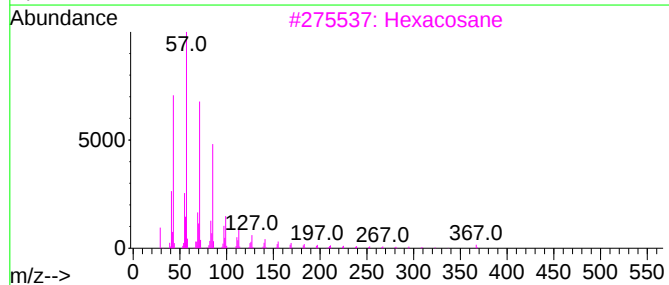
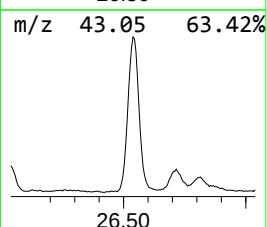
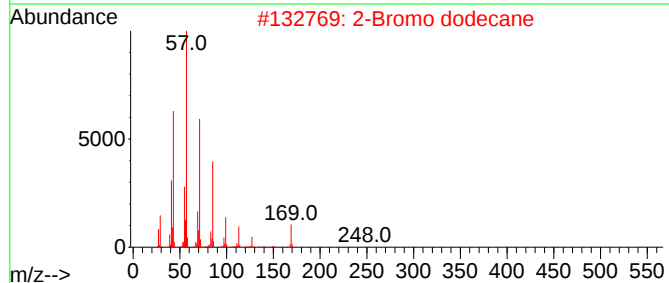
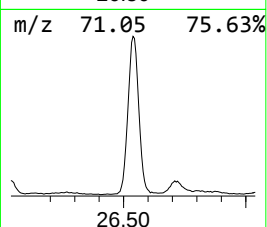
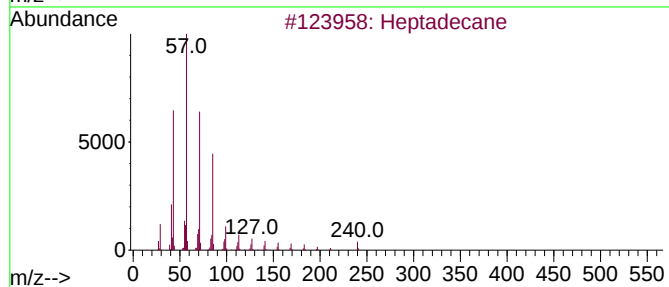
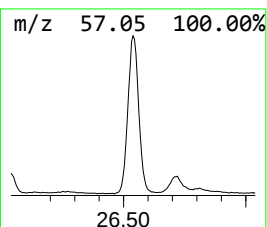
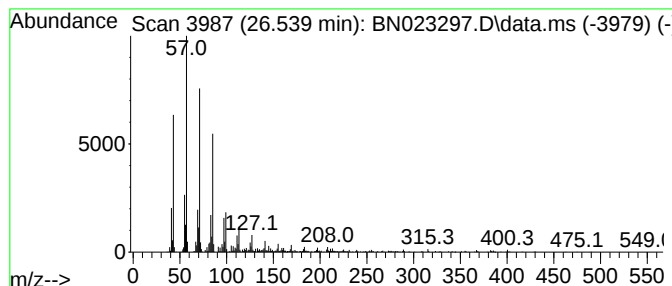
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.539	11.08 ng/ul	2624240	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	93
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
Operator : CG/JU
Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBXT9

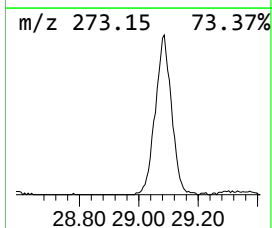
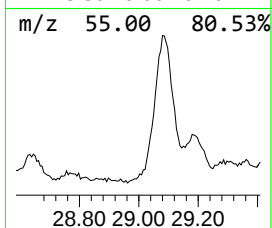
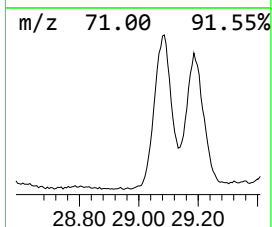
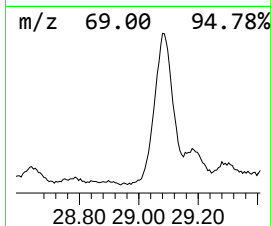
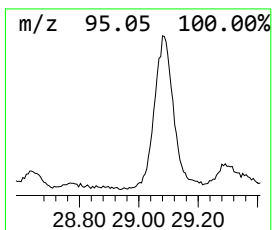
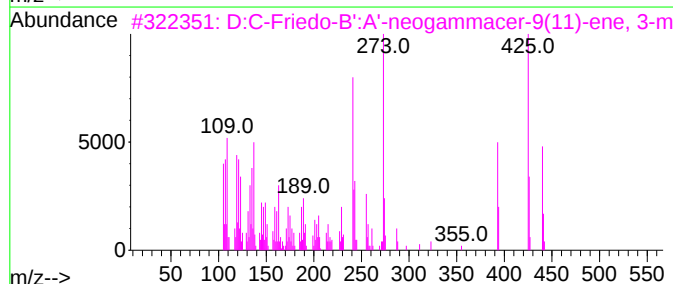
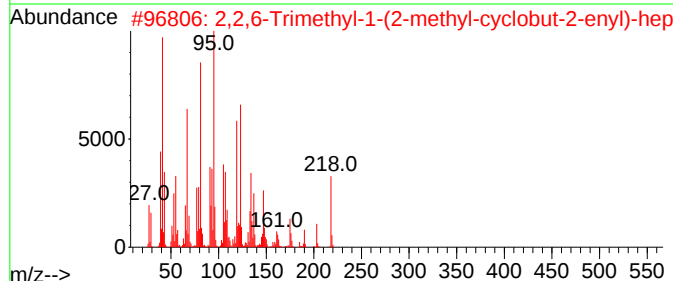
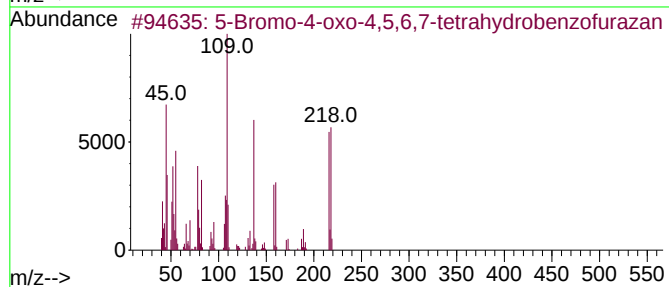
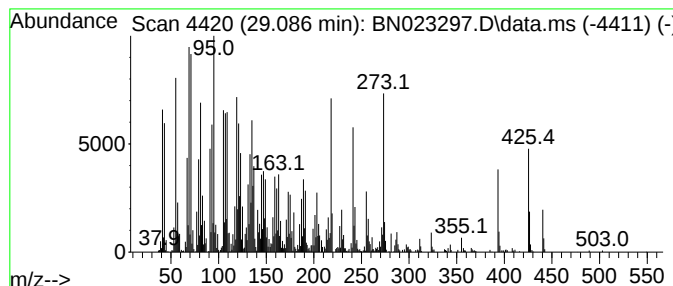
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.086	12.30 ng/ul	2913730	Perylene-d12	24.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	91
2			2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	42
3			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	38
4			6-Nitro-2-phenylthieno[2,3-d]pyr...	273	C12H7N3O3S	1000460-07-6	25
5			Sesquirosefuran	218	C15H22O	039007-93-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122022\
Data File : BN023297.D
Acq On : 20 Dec 2022 14:14
Operator : CG/JU
Sample : N6003-19
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBXT9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN121522.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
2-Pentanone, 4-...	5.063	2.3	ng/ul	202631	1	7.998	1729020	20.0
Benzophenone	16.004	10.3	ng/ul	1990370	3	14.645	3874610	20.0
n-Hexadecanoic ...	18.286	7.5	ng/ul	1789060	4	17.392	4744510	20.0
Isopropyl palmi...	18.710	2.1	ng/ul	504271	4	17.392	4744510	20.0
Octadecanoic acid	19.563	2.8	ng/ul	830585	5	21.580	5999720	20.0
1-Heneicosanol	21.274	19.6	ng/ul	5864310	5	21.580	5999720	20.0
(DEL) Alkane: S...	23.262	3.3	ng/ul	771608	6	24.027	4738530	20.0
(DEL) Alkane: S...	24.651	7.3	ng/ul	1736610	6	24.027	4738530	20.0
1-Heptacosanol	24.750	3.3	ng/ul	791542	6	24.027	4738530	20.0
(DEL) Alkane: S...	26.539	11.1	ng/ul	2624240	6	24.027	4738530	20.0
5-Bromo-4-oxo-4...	29.086	12.3	ng/ul	2913730	6	24.027	4738530	20.0