

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006486.D
 Acq On : 22 Jun 2019 13:30
 Operator : HP/JU
 Sample : K3362-14
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4203

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M

Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.052	8	11	18	rVB	25183	43688	0.73%	0.045%
2	3.134	22	25	31	rVB	62683	82175	1.36%	0.085%
3	3.646	108	112	120	rVV	79800	108758	1.81%	0.112%
4	3.764	127	132	139	rVB	44106	71326	1.18%	0.073%
5	4.287	216	221	230	rVB	358637	487141	8.09%	0.501%
6	4.752	294	300	311	rVV	37225	60604	1.01%	0.062%
7	4.875	314	321	328	rVB3	39842	67478	1.12%	0.069%
8	6.311	559	565	583	rVB	121858	205085	3.41%	0.211%
9	6.758	634	641	644	rBV	195997	283870	4.71%	0.292%
10	6.799	644	648	664	rVV	1168228	1934336	32.12%	1.990%
11	6.964	666	676	686	rVV	996336	1576850	26.19%	1.622%
12	7.058	686	692	702	rVB	194010	299905	4.98%	0.309%
13	7.158	702	709	721	rBV	1245016	1981174	32.90%	2.038%
14	7.622	780	788	797	rBV	1251908	2014902	33.46%	2.073%
15	8.334	903	909	925	rVV	1242149	2082469	34.58%	2.143%
16	8.775	978	984	993	rBV	1142683	1931416	32.08%	1.987%
17	9.187	1046	1054	1060	rVB	134556	189397	3.15%	0.195%
18	9.493	1099	1106	1123	rBV	954441	1654270	27.47%	1.702%
19	10.034	1191	1198	1214	rBV	1451321	2553853	42.41%	2.628%
20	10.405	1253	1261	1273	rBV	1849758	3028865	50.30%	3.116%
21	10.546	1278	1285	1307	rVB	971887	1811436	30.08%	1.864%
22	11.681	1470	1478	1486	rBV	86151	127250	2.11%	0.131%
23	12.004	1528	1533	1541	rVB2	23542	40490	0.67%	0.042%
24	13.169	1723	1731	1736	rBV	70793	114591	1.90%	0.118%
25	13.593	1797	1803	1811	rVB3	55568	77987	1.30%	0.080%
26	13.693	1813	1820	1824	rVV	2752382	3790408	62.95%	3.900%
27	13.740	1824	1828	1837	rVB	786678	1068117	17.74%	1.099%
28	13.969	1860	1867	1880	rBV	3010500	4506748	74.84%	4.637%
29	14.281	1913	1920	1927	rBV2	2643708	3898360	64.74%	4.011%
30	14.345	1927	1931	1939	rVB3	52941	83813	1.39%	0.086%
31	14.504	1952	1958	1977	rBV	770470	1468153	24.38%	1.511%
32	14.657	1980	1984	1991	rVV5	26456	65619	1.09%	0.068%
33	15.087	2051	2057	2063	rBV	91390	123274	2.05%	0.127%
34	15.281	2083	2090	2096	rBV	3891754	5474953	90.92%	5.633%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
 Title : SVOA CALIBRATION

35	15.334	2096	2099	2108	rVV2	46849	76674	1.27%	0.079%
36	15.422	2108	2114	2125	rVV	937104	1342214	22.29%	1.381%
37	15.516	2126	2130	2139	rVV	52111	97692	1.62%	0.101%
38	15.763	2165	2172	2179	rVV6	32504	63134	1.05%	0.065%
39	16.522	2294	2301	2305	rVB	42510	59026	0.98%	0.061%
40	16.792	2344	2347	2350	rVV3	30182	43357	0.72%	0.045%
41	16.851	2354	2357	2363	rVB	31018	45707	0.76%	0.047%
42	16.945	2368	2373	2379	rBV2	54885	91503	1.52%	0.094%
43	17.034	2379	2388	2392	rVV2	3133528	4434236	73.64%	4.562%
44	17.075	2392	2395	2400	rVV	795035	1083013	17.99%	1.114%
45	17.134	2400	2405	2417	rVV2	3749769	5449515	90.50%	5.607%
46	17.445	2450	2458	2465	rBV2	48399	103771	1.72%	0.107%
47	17.587	2479	2482	2487	rVB	33461	40574	0.67%	0.042%
48	17.822	2516	2522	2528	rBV3	64209	120557	2.00%	0.124%
49	17.881	2528	2532	2534	rBV4	32316	42591	0.71%	0.044%
50	17.916	2534	2538	2539	rVV	59337	79866	1.33%	0.082%
51	17.939	2539	2542	2550	rVV2	586234	1000013	16.61%	1.029%
52	18.004	2550	2553	2557	rVB	89739	122203	2.03%	0.126%
53	18.110	2566	2571	2575	rBV2	153834	230966	3.84%	0.238%
54	18.145	2575	2577	2583	rVB2	52095	65662	1.09%	0.068%
55	18.210	2583	2588	2596	rBV2	182254	302671	5.03%	0.311%
56	18.422	2620	2624	2628	rVB2	89273	122579	2.04%	0.126%
57	18.469	2628	2632	2637	rVB	63245	81832	1.36%	0.084%
58	18.539	2640	2644	2651	rBV2	44522	66132	1.10%	0.068%
59	18.834	2687	2694	2698	rBV4	71875	187140	3.11%	0.193%
60	18.886	2698	2703	2709	rVB3	90044	170865	2.84%	0.176%
61	18.986	2714	2720	2727	rVB3	37987	92048	1.53%	0.095%
62	19.104	2732	2740	2750	rBV	1956604	2877622	47.79%	2.961%
63	19.233	2759	2762	2770	rBV2	125038	245861	4.08%	0.253%
64	19.445	2792	2798	2801	rBV	4289762	6021469	100.00%	6.195%
65	19.475	2801	2803	2812	rVV	1535284	1786562	29.67%	1.838%
66	19.545	2812	2815	2821	rVB2	59861	92584	1.54%	0.095%
67	19.657	2830	2834	2837	rBV	76964	114819	1.91%	0.118%
68	19.798	2853	2858	2860	rBV2	78203	125831	2.09%	0.129%
69	19.939	2878	2882	2884	rBV3	69738	106762	1.77%	0.110%
70	19.980	2885	2889	2898	rVV3	268250	643118	10.68%	0.662%
71	20.080	2898	2906	2909	rVV3	164767	261767	4.35%	0.269%

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72	20.133	2909	2915	2919	rVV2	119203	226935	3.77%	0.233%
73	20.169	2919	2921	2926	rVB	69317	87915	1.46%	0.090%
74	20.275	2937	2939	2943	rVV	34469	37752	0.63%	0.039%
75	20.328	2945	2948	2951	rVB2	55581	70503	1.17%	0.073%
76	20.422	2961	2964	2967	rBV4	60754	79700	1.32%	0.082%
77	20.504	2973	2978	2982	rBV	886306	982500	16.32%	1.011%
78	20.545	2982	2985	2988	rVB	108452	108978	1.81%	0.112%
79	20.639	2999	3001	3005	rVB2	47186	50826	0.84%	0.052%
80	20.686	3006	3009	3013	rVV3	61324	87646	1.46%	0.090%
81	20.733	3014	3017	3025	rVB	93769	123089	2.04%	0.127%
82	20.898	3041	3045	3048	rBV2	144312	200066	3.32%	0.206%
83	20.933	3048	3051	3054	rVV2	109070	138120	2.29%	0.142%
84	20.975	3054	3058	3065	rVV3	472227	823191	13.67%	0.847%
85	21.169	3087	3091	3096	rVB	1917531	2025742	33.64%	2.084%
86	21.257	3098	3106	3110	rVV2	2902108	4781002	79.40%	4.919%
87	21.292	3110	3112	3122	rVB	823088	950900	15.79%	0.978%
88	21.410	3128	3132	3136	rBV	369349	421170	6.99%	0.433%
89	21.804	3197	3199	3202	rVB2	76794	81457	1.35%	0.084%
90	21.845	3202	3206	3209	rBV	367002	443731	7.37%	0.457%
91	22.304	3279	3284	3289	rBV	467742	629619	10.46%	0.648%
92	22.804	3364	3369	3377	rBV2	881272	2354363	39.10%	2.422%
93	23.298	3448	3453	3456	rBV	344522	578467	9.61%	0.595%
94	23.351	3456	3462	3467	rBV	2276274	4226937	70.20%	4.349%
95	23.492	3479	3486	3491	rBV2	1691486	3099779	51.48%	3.189%
96	23.998	3566	3572	3581	rVB	548548	1083878	18.00%	1.115%
97	24.092	3583	3588	3597	rBV2	124232	307916	5.11%	0.317%
98	24.727	3690	3696	3703	rBV	293672	672064	11.16%	0.691%
99	25.580	3836	3841	3856	rVB	215928	558914	9.28%	0.575%
100	25.721	3859	3865	3876	rVB	273865	757593	12.58%	0.779%

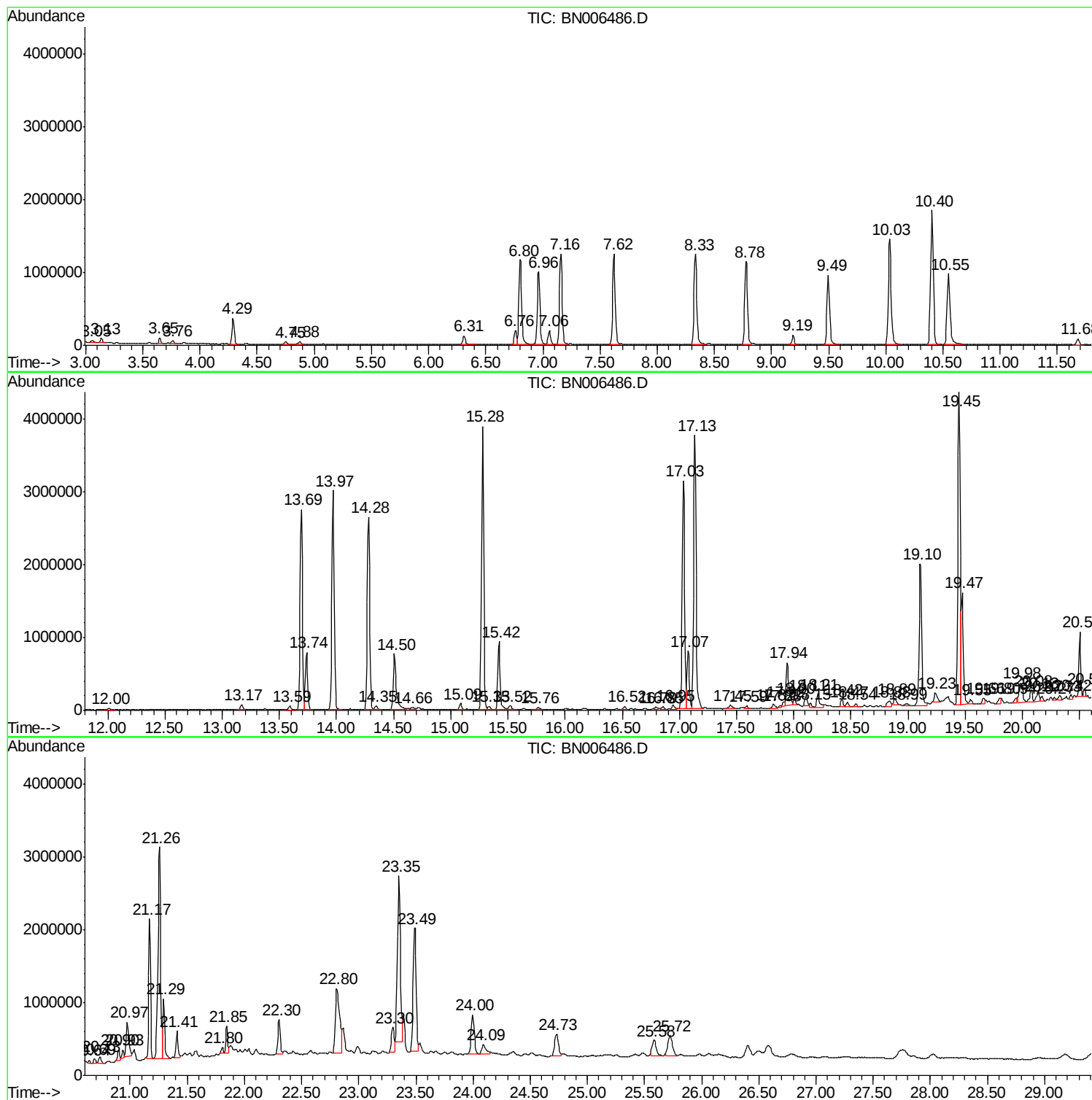
Sum of corrected areas: 97191450

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

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



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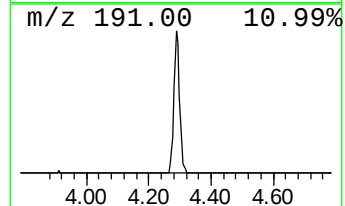
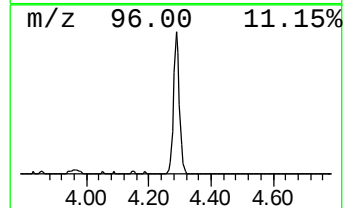
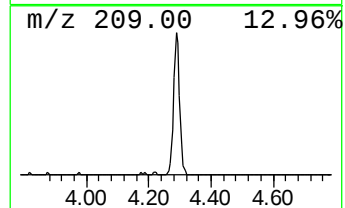
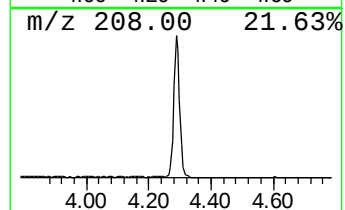
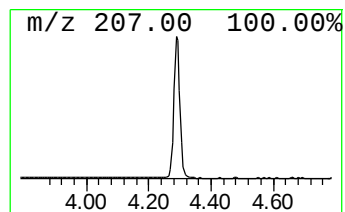
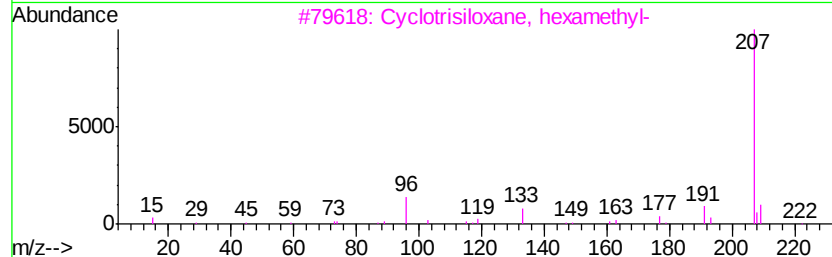
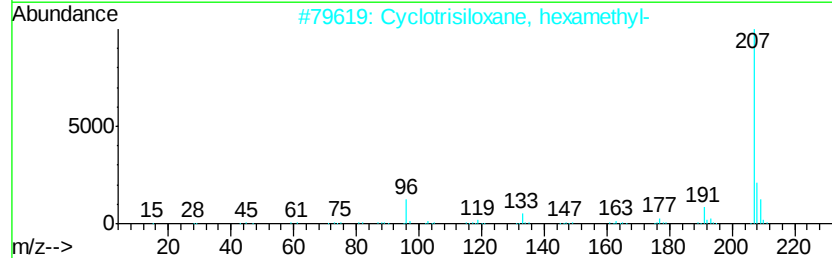
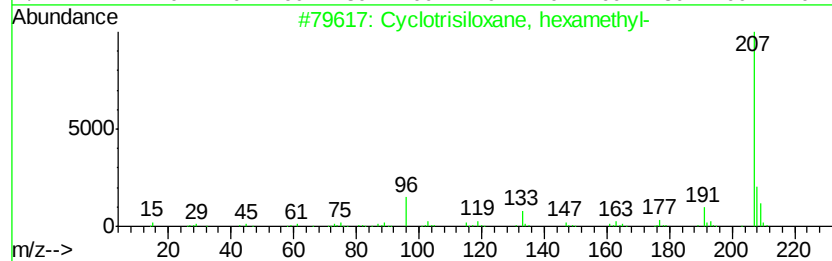
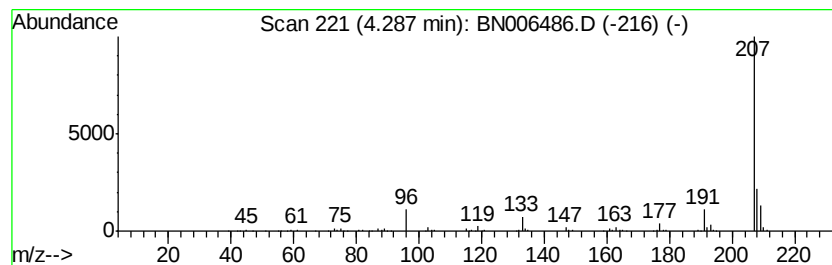
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclotrisiloxane, hexamethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.29	4.84 ng/ul	487141	1,4-Dichlorobenzene-d4	7.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	91
2			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	91
3			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	80
4			Arsenous acid, tris(trimethylsil...	342	C9H27AsO3Si3	055429-29-3	78
5			4-Cyclohexene-1,2-dicarboximide,...	207	C12H17N02	028916-00-9	42



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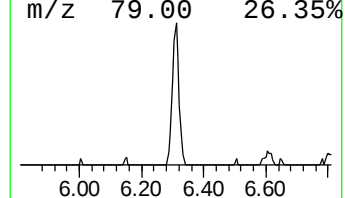
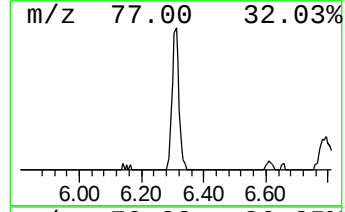
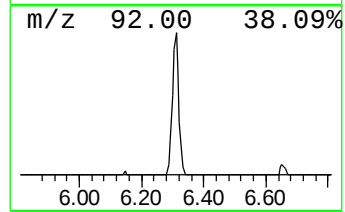
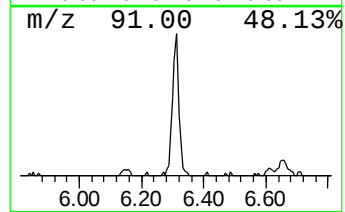
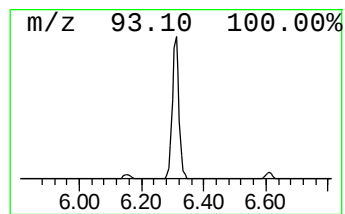
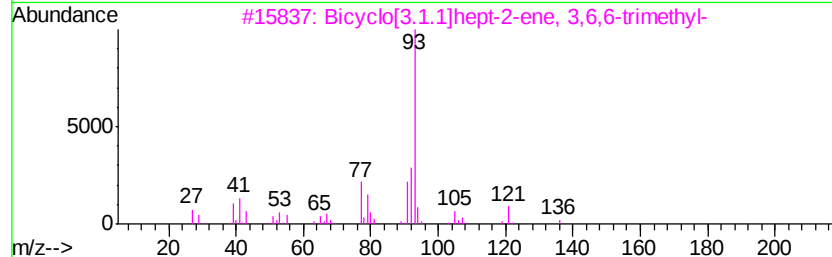
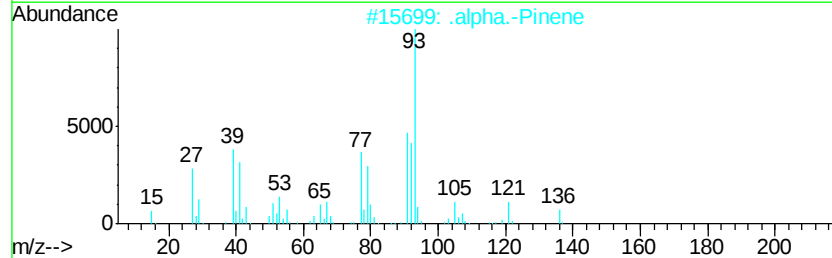
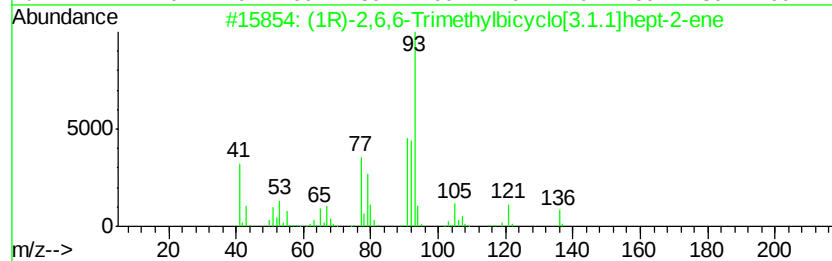
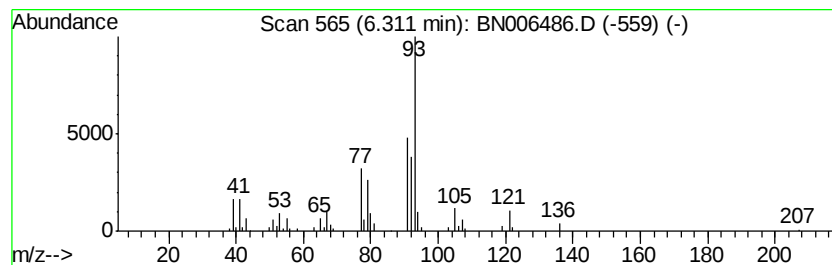
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	2.04 ng/ul	205085	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene	136	C10H16	007785-70-8	95
2		.alpha.-Pinene	136	C10H16	000080-56-8	94
3		Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-	136	C10H16	004889-83-2	94
4		.alpha.-Pinene	136	C10H16	000080-56-8	94
5		Tricyclo[2.2.1.0(2,6)]heptane, 1,4-dichloro-	136	C10H16	000508-32-7	91



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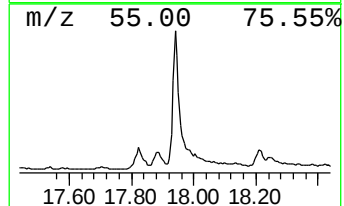
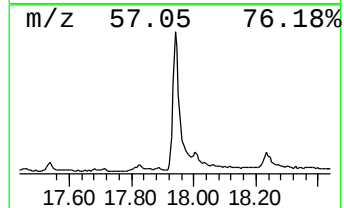
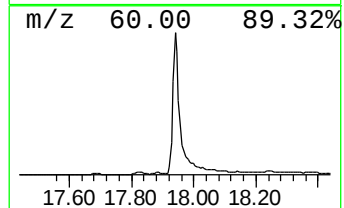
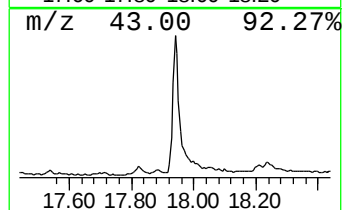
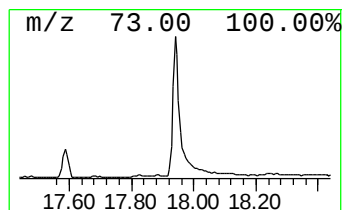
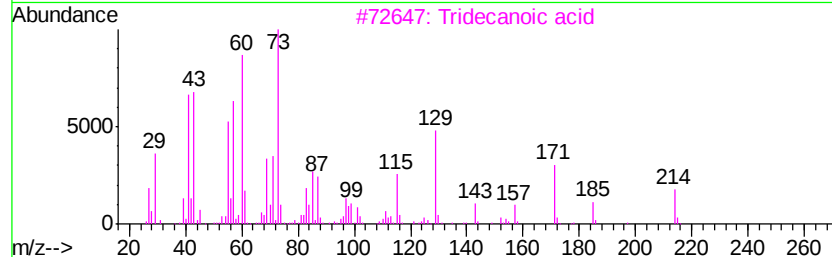
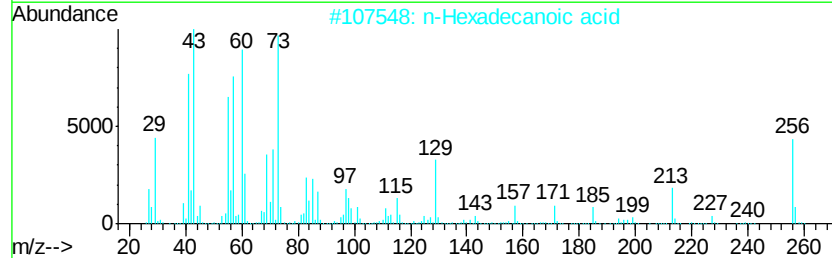
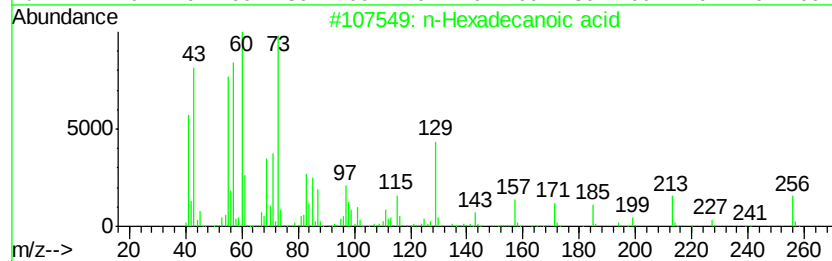
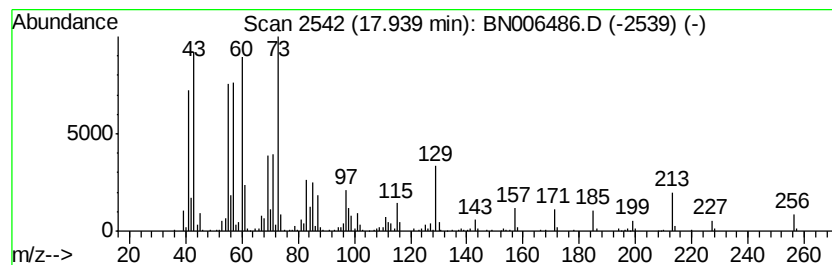
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.94	4.51 ng/ul	1000010	Phenanthrene-d10	17.03		
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	95	
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	93	
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90	



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Data File : BN006486.D
Acq On : 22 Jun 2019 13:30
Operator : HP/JU
Sample : K3362-14
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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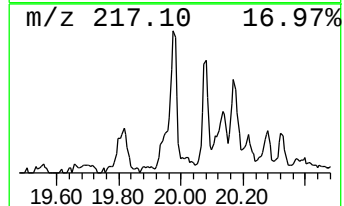
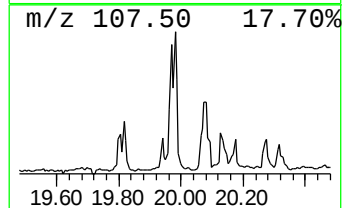
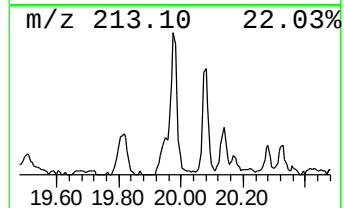
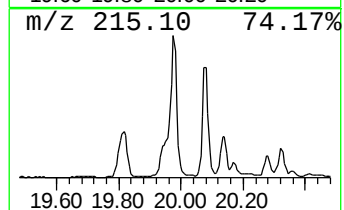
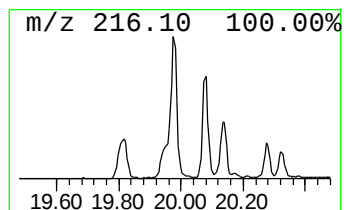
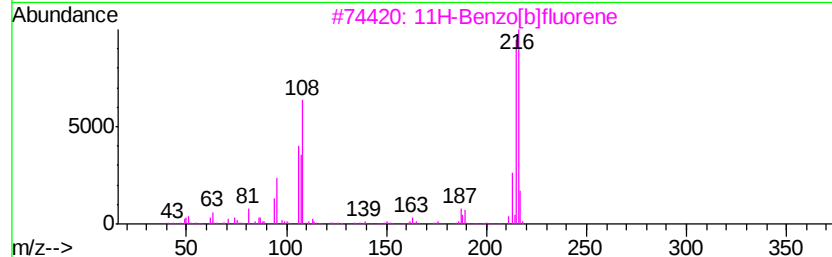
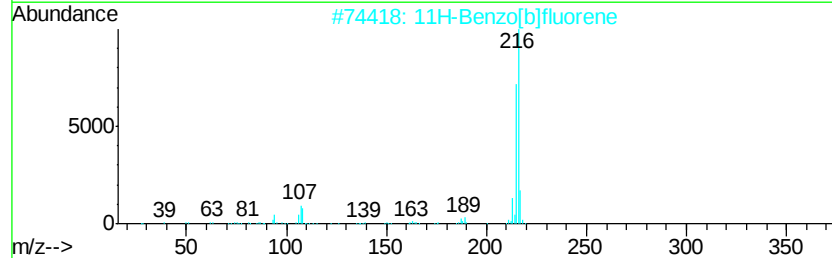
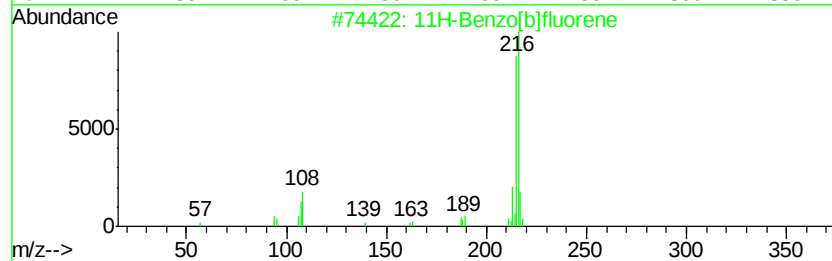
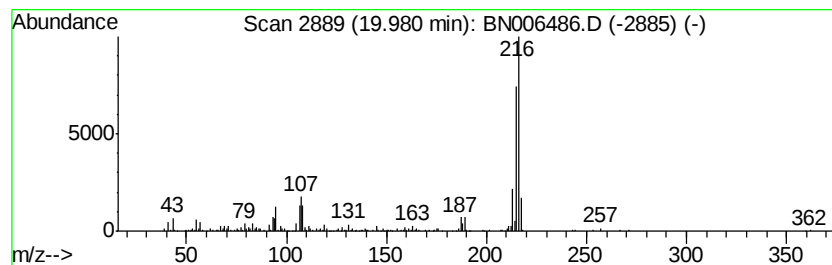
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.98	2.69 ng/ul	643118	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
Data File : BN006486.D
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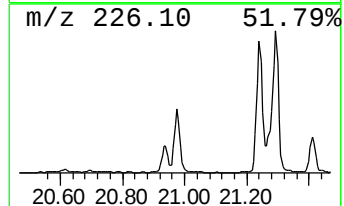
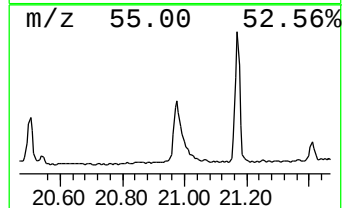
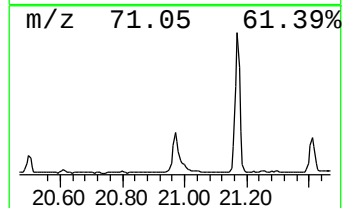
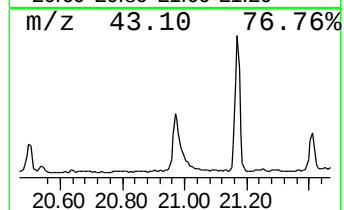
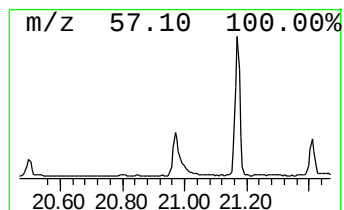
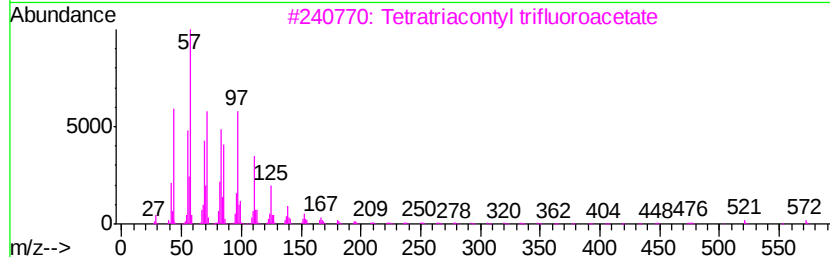
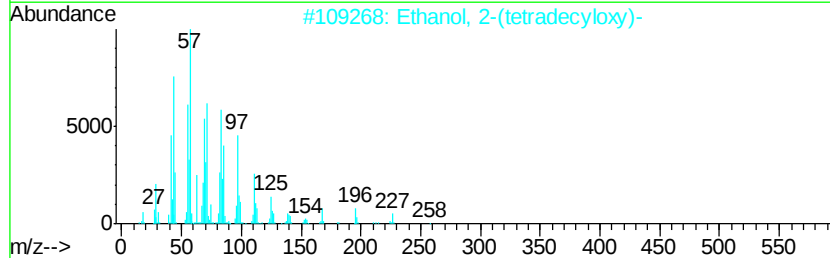
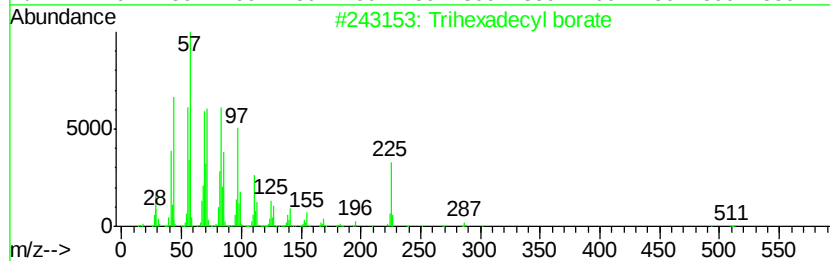
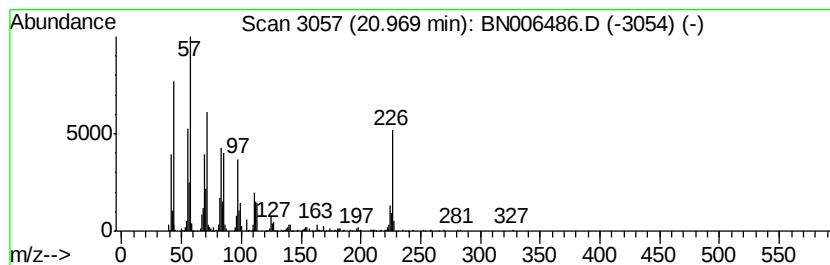
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Trihexadecyl borate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	3.44 ng/ul	823191	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trihexadecyl borate	735	C48H99B03	002665-11-4	53
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	50
3		Tetratriacontyl trifluoroacetate	591	C36H69F3O2	1000351-75-3	49
4		Tetratriacontyl pentafluoropropi...	641	C37H69F5O2	1000351-81-5	49
5		Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
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Sample : K3362-14
Misc :
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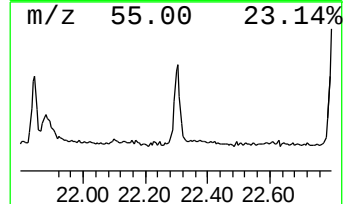
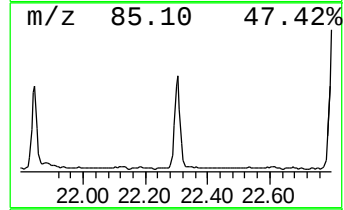
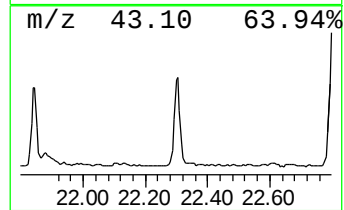
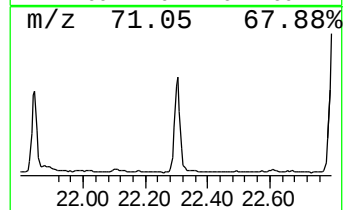
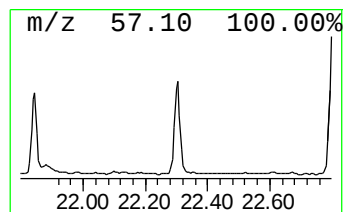
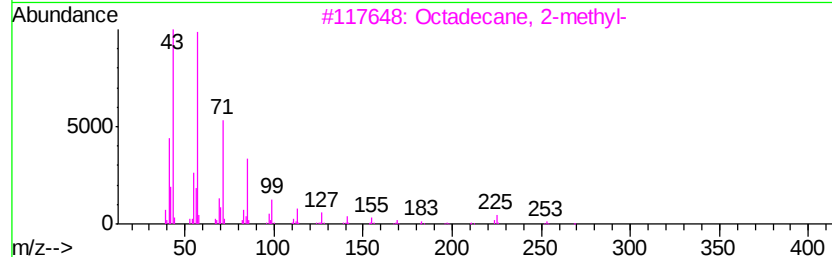
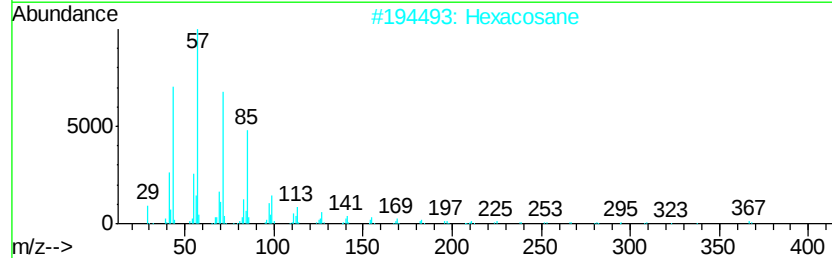
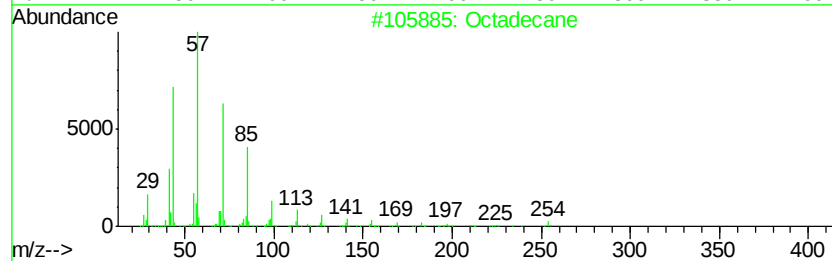
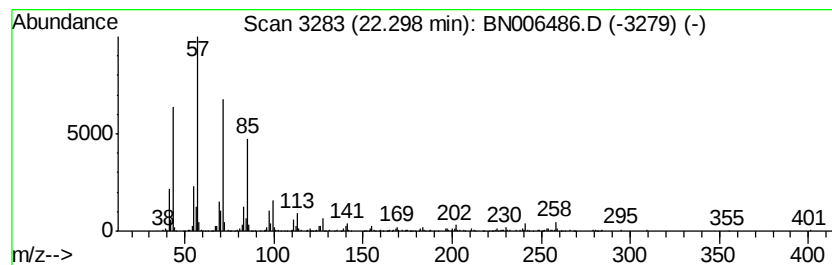
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.30	2.63 ng/ul	629619	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	96
2		Hexacosane	366	C26H54	000630-01-3	93
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
4		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
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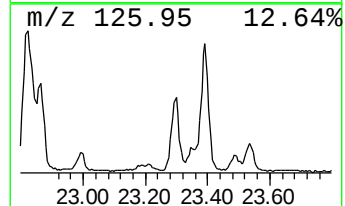
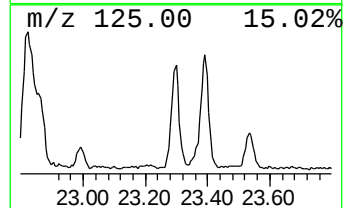
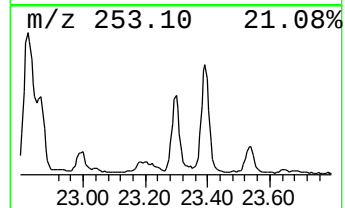
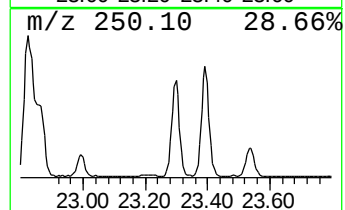
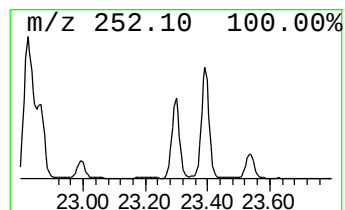
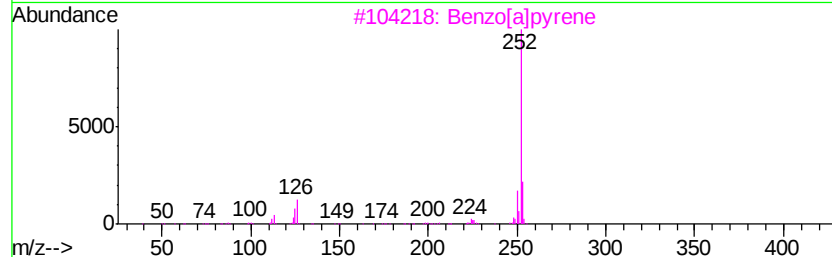
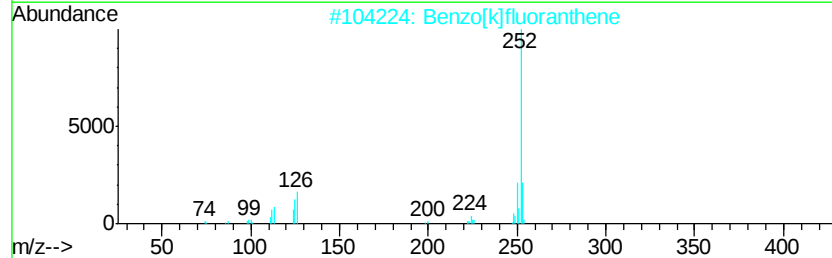
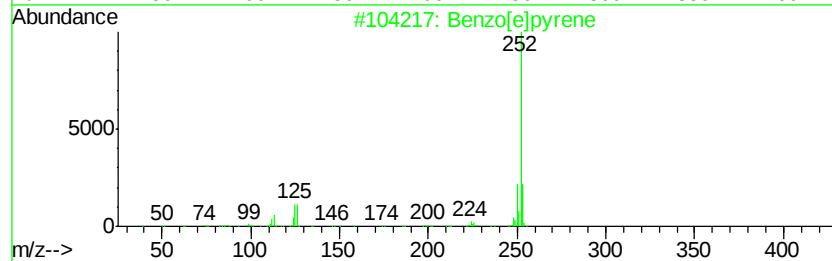
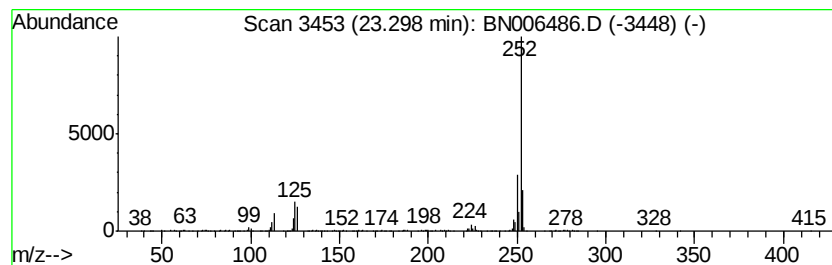
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.30	3.73 ng/ul	578467	Perylene-d12	23.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Benzo[a]pyrene	252	C20H12	000050-32-8	95
4		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94
5		Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
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Acq On : 22 Jun 2019 13:30
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Sample : K3362-14
Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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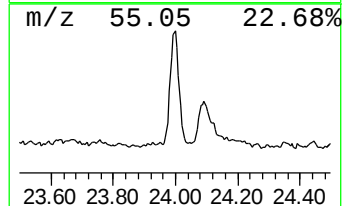
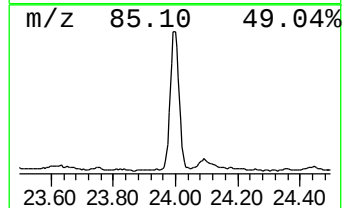
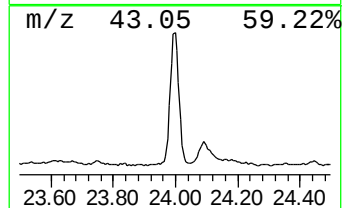
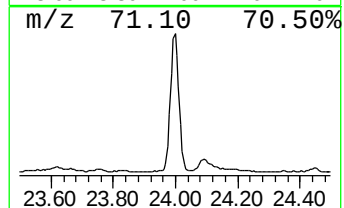
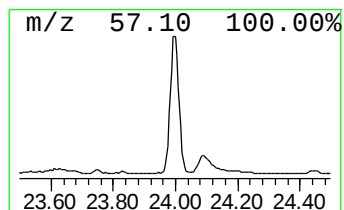
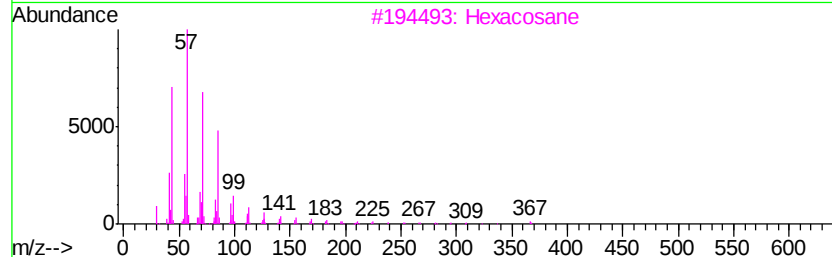
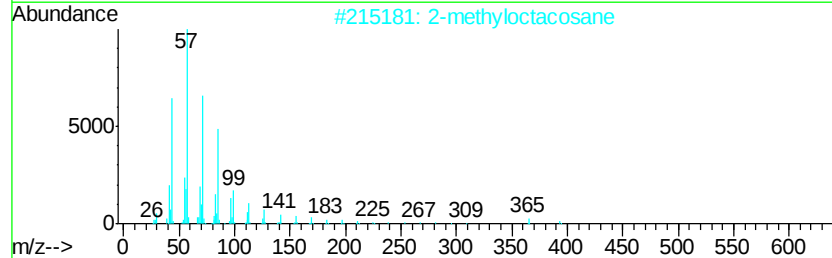
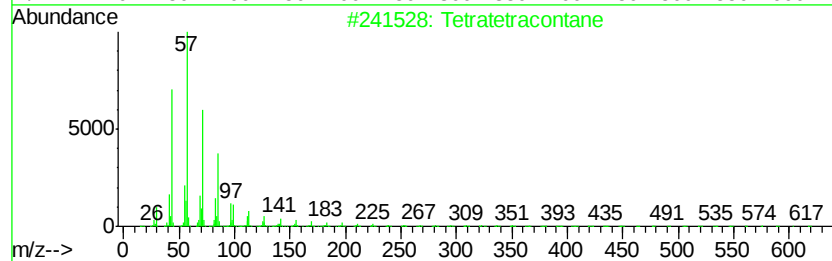
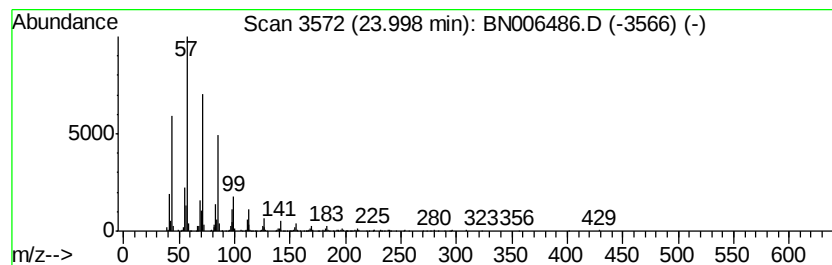
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.00	6.99 ng/ul	1083880	Perylene-d12	23.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
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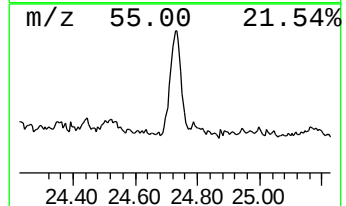
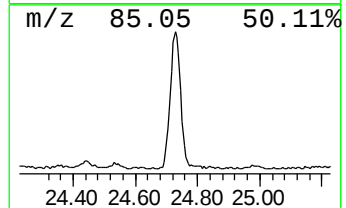
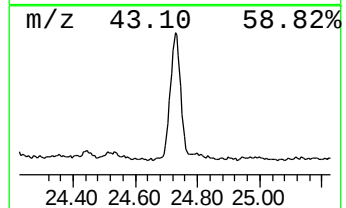
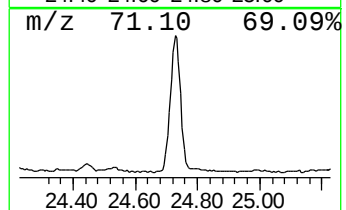
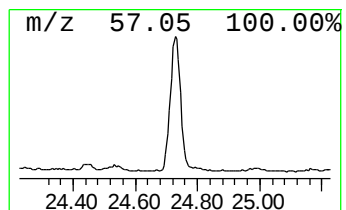
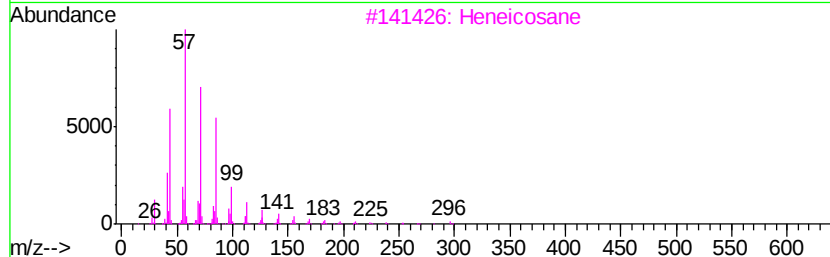
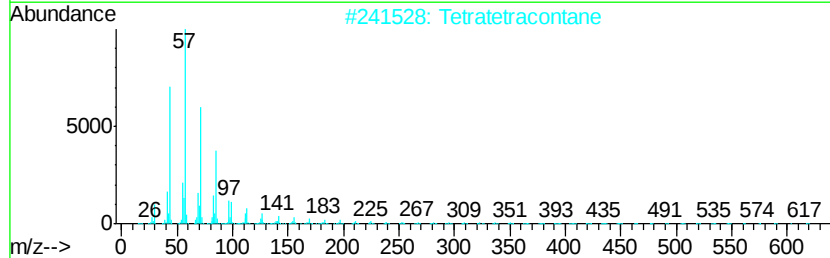
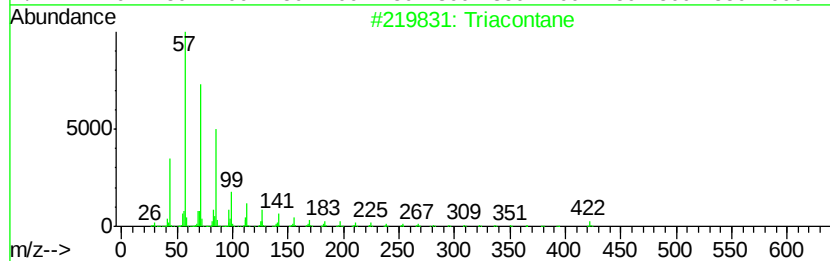
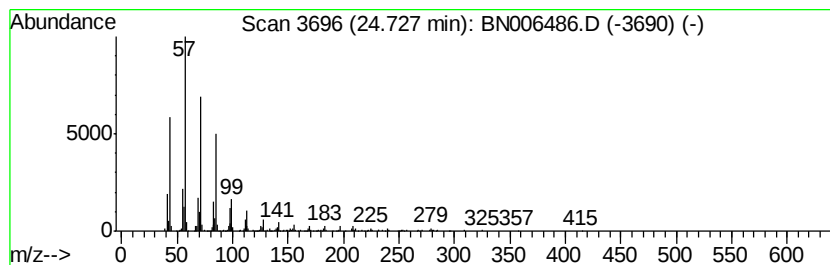
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.73	4.34 ng/ul	672064	Perylene-d12	23.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontane	422	C30H62	000638-68-6	91
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Hexacosane	366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
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Sample : K3362-14
Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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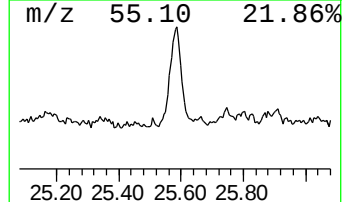
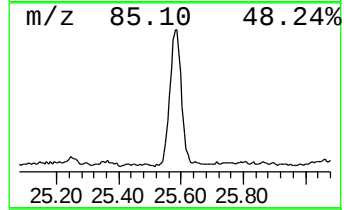
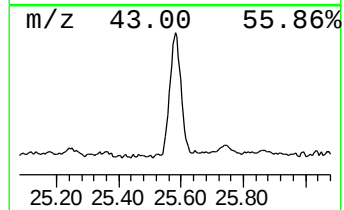
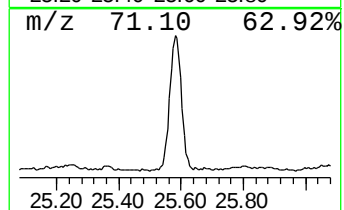
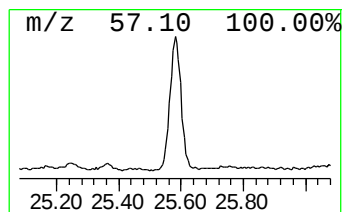
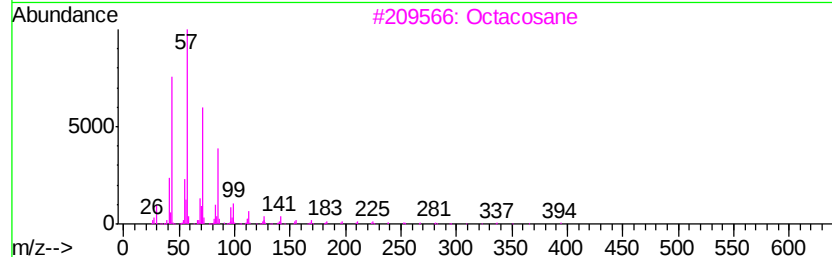
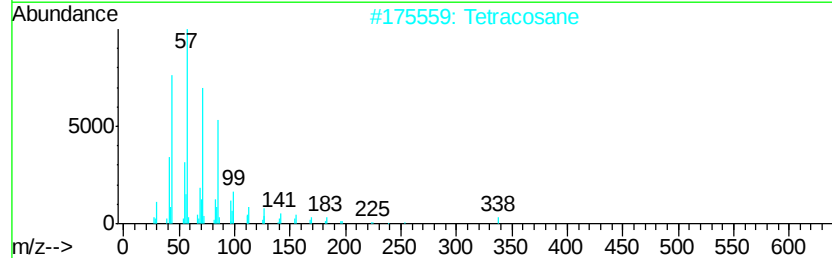
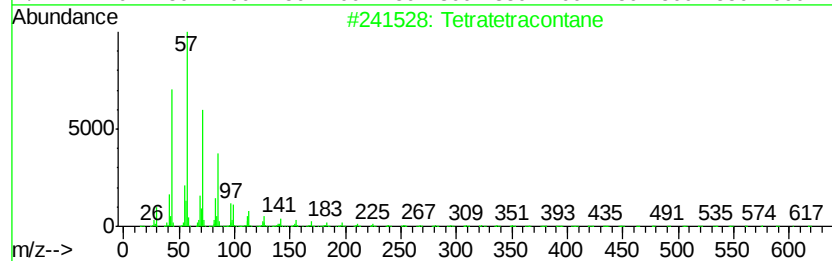
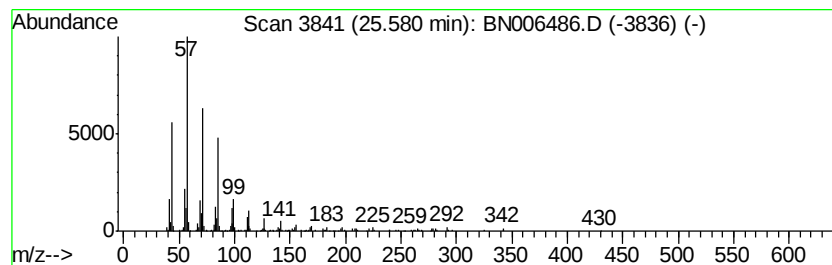
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.58	3.61 ng/ul	558914	Perylene-d12	23.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	90
2			Tetracosane	338	C24H50	000646-31-1	87
3			Octacosane	394	C28H58	000630-02-4	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006486.D
 Acq On : 22 Jun 2019 13:30
 Operator : HP/JU
 Sample : K3362-14
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4203

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061919MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclotrisiloxane,...	4.29	4.8	ng/ul	487141	1	7.62	2014900	20.0
(1R)-2,6,6-Trimet...	6.31	2.0	ng/ul	205085	1	7.62	2014900	20.0
n-Hexadecanoic acid	17.94	4.5	ng/ul	1000010	4	17.03	4434240	20.0
11H-Benzo[b]fluorene	19.98	2.7	ng/ul	643118	5	21.26	4781000	20.0
unknown-01	20.50	4.1	ng/ul	982500	5	21.26	4781000	20.0
Trihexadecyl borate	20.97	3.4	ng/ul	823191	5	21.26	4781000	20.0
(DEL) Alkane: Str...	22.30	2.6	ng/ul	629619	5	21.26	4781000	20.0
Benzo[e]pyrene	23.30	3.7	ng/ul	578467	6	23.49	3099780	20.0
(DEL) Alkane: Str...	24.00	7.0	ng/ul	1083880	6	23.49	3099780	20.0
(DEL) Alkane: Str...	24.73	4.3	ng/ul	672064	6	23.49	3099780	20.0
(DEL) Alkane: Str...	25.58	3.6	ng/ul	558914	6	23.49	3099780	20.0