

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
 Data File : BP001869.D
 Acq On : 24 Feb 2020 20:00
 Operator : CG/JU
 Sample : L1536-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB6X3

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_P\METHODS\SOM-EPA-BP021820MA.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	2.922	3	6	27	rVB	2400505	3476666	22.94%	2.432%
2	3.170	44	48	61	rBV	111543	157547	1.04%	0.110%
3	3.811	153	157	163	rVB	31732	52110	0.34%	0.036%
4	3.899	168	172	178	rVB	64718	90382	0.60%	0.063%
5	4.122	206	210	215	rBV	41780	54970	0.36%	0.038%
6	4.258	229	233	238	rVB2	47494	69108	0.46%	0.048%
7	4.799	320	325	334	rVB	42899	68212	0.45%	0.048%
8	5.734	478	484	490	rVB4	29242	61056	0.40%	0.043%
9	6.369	585	592	597	rBV	46387	72949	0.48%	0.051%
10	6.828	657	670	685	rBV	1553933	2537911	16.75%	1.775%
11	6.993	691	698	707	rBV	1271676	1967165	12.98%	1.376%
12	7.187	724	731	742	rBV	1732352	2733377	18.04%	1.912%
13	7.358	755	760	766	rVB	34752	50747	0.33%	0.035%
14	7.652	803	810	818	rBV	1269879	1969350	13.00%	1.378%
15	7.728	818	823	830	rVB	29670	44290	0.29%	0.031%
16	7.887	845	850	856	rBV5	21015	39722	0.26%	0.028%
17	8.357	918	930	942	rVV	2038731	3272793	21.60%	2.289%
18	8.793	992	1004	1016	rBV	1562396	2587954	17.08%	1.810%
19	8.893	1016	1021	1029	rVB4	19289	43465	0.29%	0.030%
20	8.975	1029	1035	1043	rBV	171869	261890	1.73%	0.183%
21	9.146	1056	1064	1069	rBV	70872	116711	0.77%	0.082%
22	9.287	1082	1088	1097	rBV	112711	182850	1.21%	0.128%
23	9.510	1119	1126	1134	rVV	1280738	2096020	13.83%	1.466%
24	10.052	1210	1218	1230	rVV	2134511	3611163	23.83%	2.526%
25	10.422	1274	1281	1288	rBV	1709908	2957922	19.52%	2.069%
26	10.475	1288	1290	1296	rVB	28353	39181	0.26%	0.027%
27	10.557	1296	1304	1321	rBV	1851639	3282633	21.66%	2.296%
28	12.016	1547	1552	1560	rVB2	60398	107731	0.71%	0.075%
29	12.093	1560	1565	1575	rVB2	21214	43252	0.29%	0.030%
30	13.704	1828	1839	1844	rBV	3731291	5262364	34.73%	3.681%
31	13.751	1844	1847	1851	rVB	189111	239413	1.58%	0.167%
32	13.975	1876	1885	1898	rBV	4349197	6760339	44.62%	4.729%
33	14.287	1931	1938	1945	rBV2	2729438	4095285	27.03%	2.865%
34	14.351	1945	1949	1953	rBV2	38340	60214	0.40%	0.042%

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
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35	14.487	1965	1972	1982	rBV	1640811	2421254	15.98%	1.694%
36	14.640	1993	1998	2003	rBV4	28673	58536	0.39%	0.041%
37	14.692	2003	2007	2017	rVB4	39452	105884	0.70%	0.074%
38	15.287	2101	2108	2114	rBV	5922807	8222828	54.27%	5.752%
39	15.339	2114	2117	2122	rVB	46509	66920	0.44%	0.047%
40	15.404	2122	2128	2137	rBV	1472172	2047507	13.51%	1.432%
41	15.528	2143	2149	2157	rVB	69357	118692	0.78%	0.083%
42	15.639	2164	2168	2180	rVB4	20951	49299	0.33%	0.034%
43	15.887	2205	2210	2216	rVB5	28813	48203	0.32%	0.034%
44	16.075	2239	2242	2249	rVB2	32715	50396	0.33%	0.035%
45	16.692	2343	2347	2350	rBV	31644	48791	0.32%	0.034%
46	16.781	2358	2362	2367	rBV5	15503	37641	0.25%	0.026%
47	16.851	2368	2374	2379	rVV3	40353	98799	0.65%	0.069%
48	17.039	2400	2406	2411	rBV	3392189	5072607	33.48%	3.548%
49	17.081	2411	2413	2417	rVV	627943	756610	4.99%	0.529%
50	17.139	2417	2423	2435	rVB2	5971069	8787773	58.00%	6.147%
51	17.445	2471	2475	2481	rBV	67204	113339	0.75%	0.079%
52	17.728	2513	2523	2529	rBV	46889	163939	1.08%	0.115%
53	17.916	2549	2555	2559	rBV2	167213	274729	1.81%	0.192%
54	17.963	2559	2563	2568	rVB	190540	250571	1.65%	0.175%
55	18.033	2570	2575	2581	rVV3	77803	131899	0.87%	0.092%
56	18.098	2582	2586	2590	rBV2	171437	272580	1.80%	0.191%
57	18.139	2590	2593	2597	rVB	114977	138649	0.92%	0.097%
58	18.251	2608	2612	2618	rVB4	161036	235710	1.56%	0.165%
59	18.404	2632	2638	2647	rBV4	108723	321140	2.12%	0.225%
60	18.616	2671	2674	2676	rBV3	44521	60199	0.40%	0.042%
61	18.651	2677	2680	2683	rVB2	72927	91435	0.60%	0.064%
62	18.804	2703	2706	2711	rVB2	135401	217721	1.44%	0.152%
63	18.886	2717	2720	2725	rVB3	232023	355607	2.35%	0.249%
64	18.939	2726	2729	2733	rBV2	75927	82471	0.54%	0.058%
65	19.028	2740	2744	2745	rBV	116538	127272	0.84%	0.089%
66	19.057	2745	2749	2755	rVB	1500785	1973890	13.03%	1.381%
67	19.275	2783	2786	2789	rBV	61072	83772	0.55%	0.059%
68	19.386	2799	2805	2816	rBV3	9434796	15152401	100.00%	10.599%
69	19.586	2836	2839	2842	rBV	103116	127711	0.84%	0.089%
70	19.739	2860	2865	2869	rBV2	105767	196032	1.29%	0.137%
71	20.051	2915	2918	2921	rBV3	121170	167953	1.11%	0.117%

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72	20.092	2922	2925	2930	rVB3	223931	288126	1.90%	0.202%
73	20.186	2938	2941	2945	rVB2	96281	120404	0.79%	0.084%
74	20.233	2946	2949	2950	rVV	87608	103608	0.68%	0.072%
75	20.251	2950	2952	2956	rVB2	169637	210270	1.39%	0.147%
76	20.433	2980	2983	2986	rBV2	155318	164733	1.09%	0.115%
77	20.622	3013	3015	3021	rVB	152853	197258	1.30%	0.138%
78	20.827	3047	3050	3052	rBV	90609	91209	0.60%	0.064%
79	20.875	3053	3058	3074	rVV5	1026265	1782034	11.76%	1.247%
80	21.074	3089	3092	3095	rVB	127403	122988	0.81%	0.086%
81	21.133	3095	3102	3106	rBV2	4548701	6409189	42.30%	4.483%
82	21.169	3106	3108	3116	rVB	638329	761324	5.02%	0.533%
83	21.251	3119	3122	3126	rVV	213556	256203	1.69%	0.179%
84	21.316	3130	3133	3137	rVB2	261380	294514	1.94%	0.206%
85	21.751	3202	3207	3218	rBV4	673954	1130119	7.46%	0.791%
86	22.063	3255	3260	3271	rBV	2686867	3942342	26.02%	2.758%
87	22.210	3281	3285	3293	rVB	425213	598617	3.95%	0.419%
88	22.433	3318	3323	3331	rVB	729255	1038014	6.85%	0.726%
89	22.686	3361	3366	3367	rBV	432045	625693	4.13%	0.438%
90	22.721	3368	3372	3375	rVV	1295643	2162972	14.27%	1.513%
91	22.757	3375	3378	3389	rVB	1635813	2498671	16.49%	1.748%
92	23.163	3443	3447	3448	rBV	205636	277078	1.83%	0.194%
93	23.210	3449	3455	3461	rVV	4704917	8649925	57.09%	6.051%
94	23.292	3465	3469	3473	rVV	419085	628607	4.15%	0.440%
95	23.351	3473	3479	3485	rVB2	2329995	4278813	28.24%	2.993%
96	23.598	3516	3521	3528	rVB2	468266	816846	5.39%	0.571%
97	23.945	3574	3580	3586	rVB	1266928	2415772	15.94%	1.690%
98	24.015	3586	3592	3606	rVB	839099	1692819	11.17%	1.184%
99	25.574	3850	3857	3869	rVB2	738029	2070019	13.66%	1.448%
100	26.427	3994	4002	4018	rVB2	732137	2329170	15.37%	1.629%

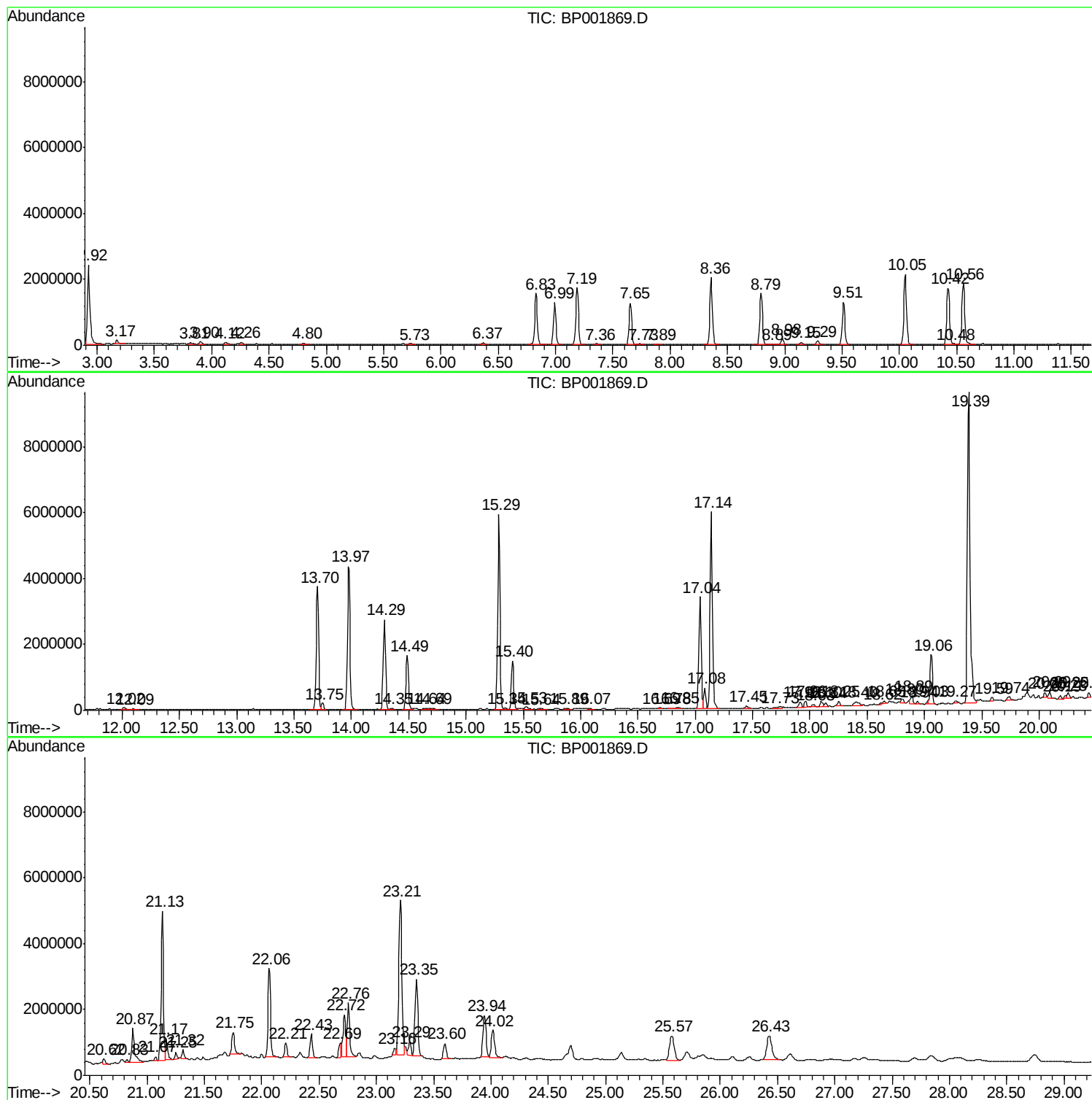
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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



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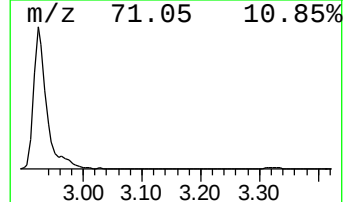
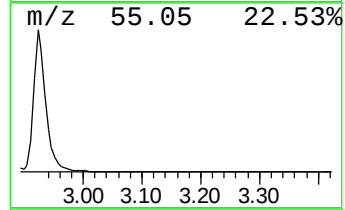
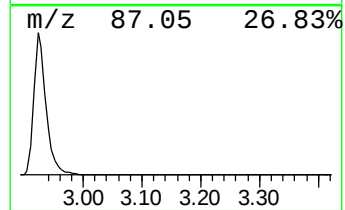
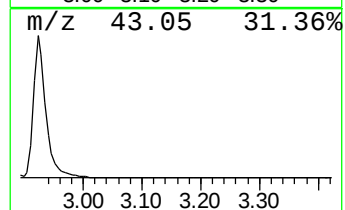
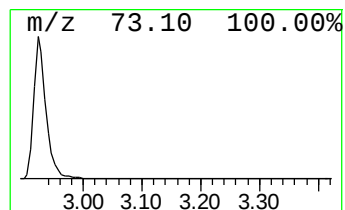
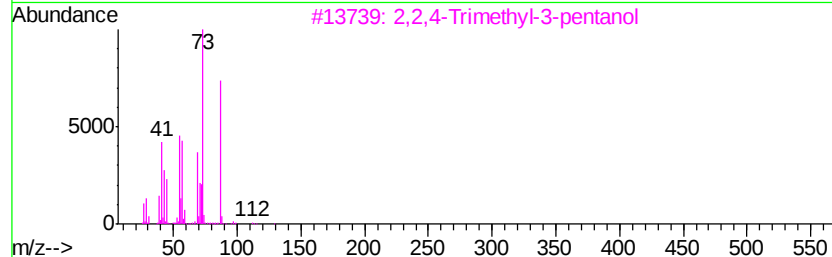
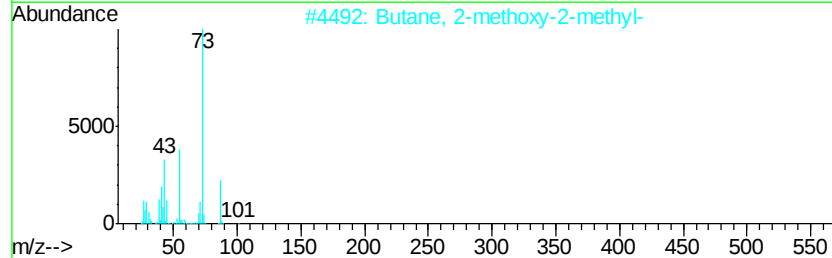
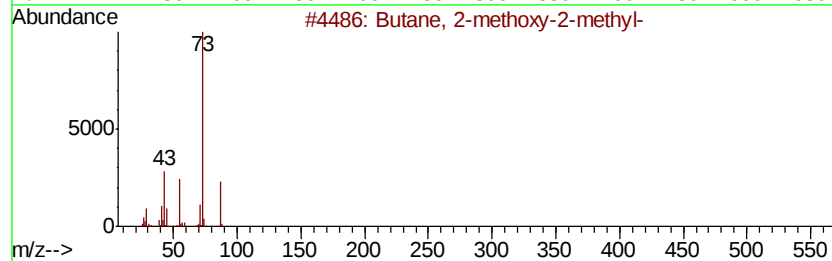
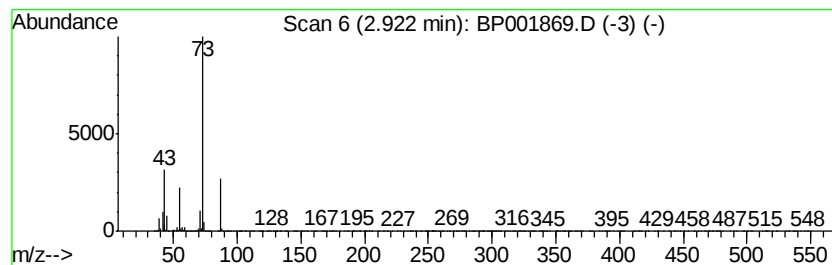
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	35.31 ng/ul	3476670	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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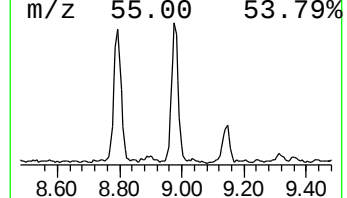
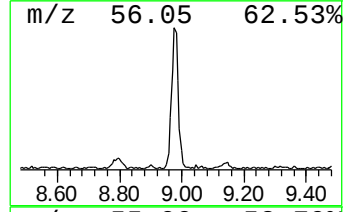
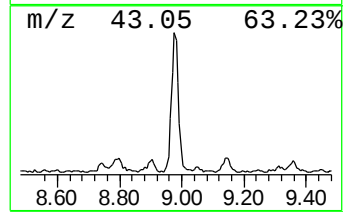
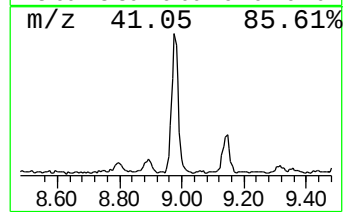
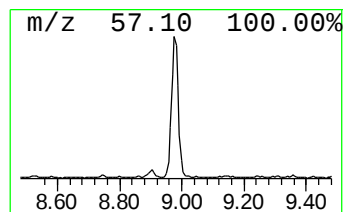
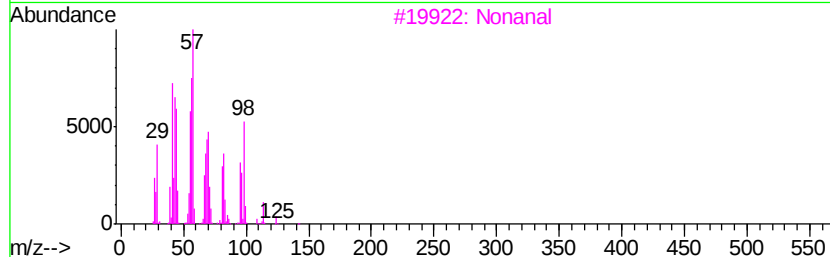
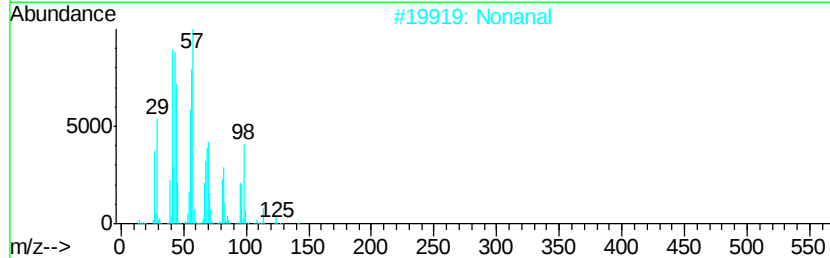
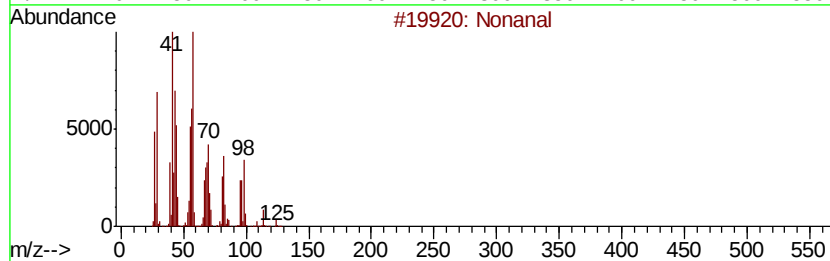
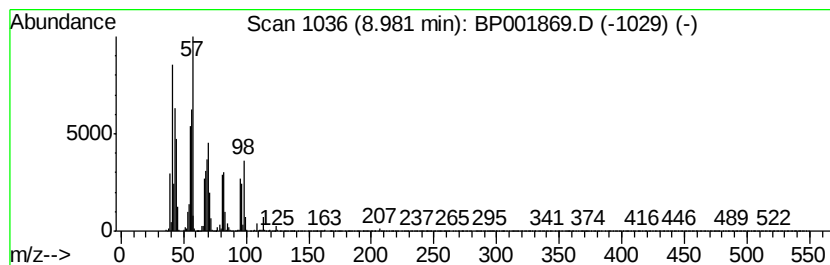
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Nonanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	2.66 ng/ul	261890	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonanal	142	C9H18O	000124-19-6	91
2		Nonanal	142	C9H18O	000124-19-6	90
3		Nonanal	142	C9H18O	000124-19-6	90
4		Nonanal	142	C9H18O	000124-19-6	72
5		Heptane, 2-chloro-	134	C7H15Cl	001001-89-4	27



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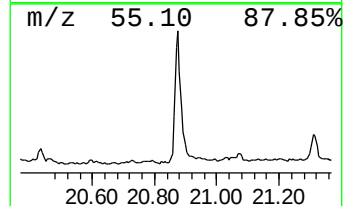
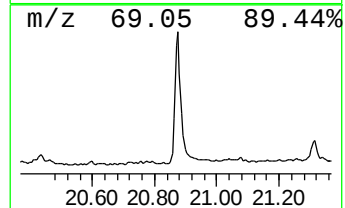
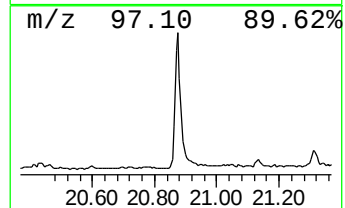
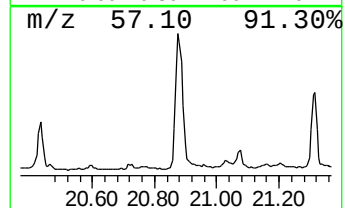
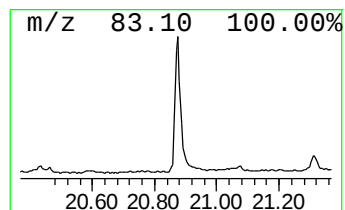
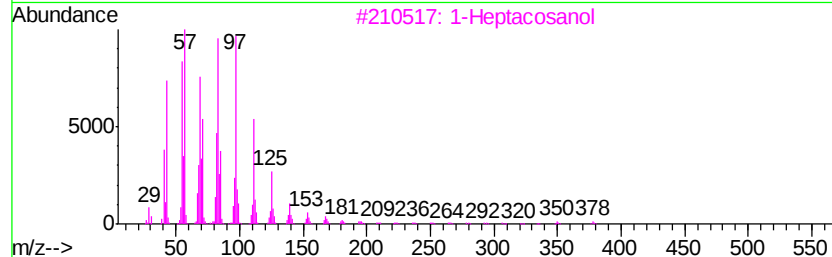
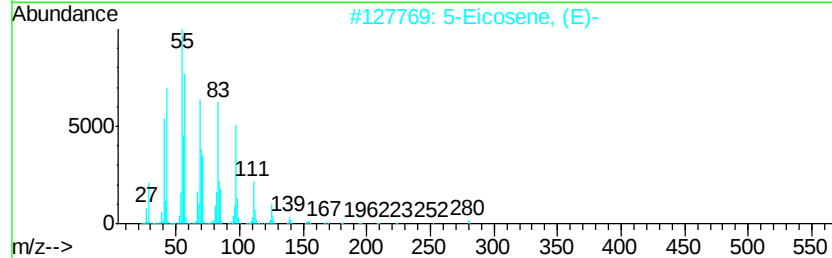
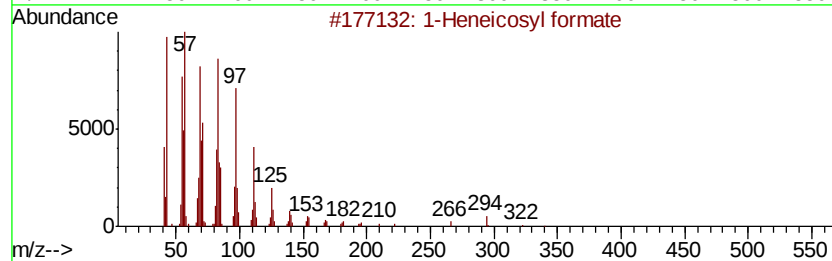
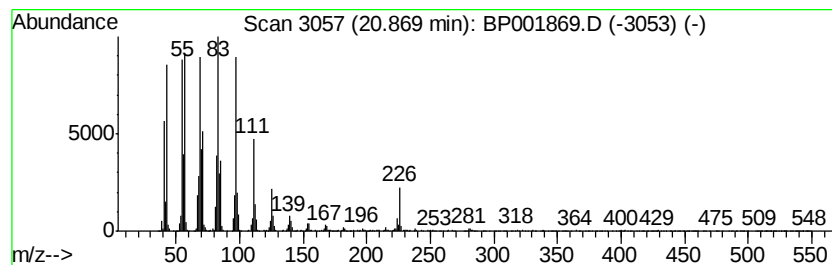
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Peak Number 3 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	5.56 ng/ul	1782030	Chrysene-d12	21.13
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Heneicosyl formate	340 C22H44O2	077899-03-7	98
2	5-Eicosene, (E)-	280 C20H40	074685-30-6	98
3	1-Heptacosanol	396 C27H56O	002004-39-9	95
4	Behenic alcohol	326 C22H46O	000661-19-8	95
5	n-Tetracosanol-1	354 C24H50O	000506-51-4	95



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Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
Operator : CG/JU
Sample : L1536-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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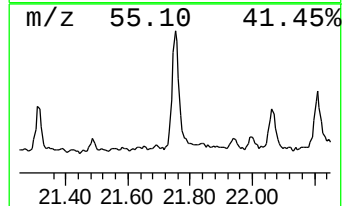
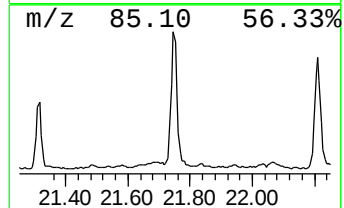
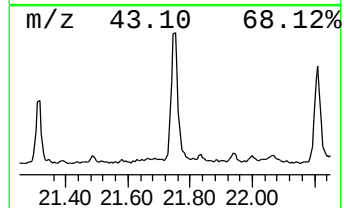
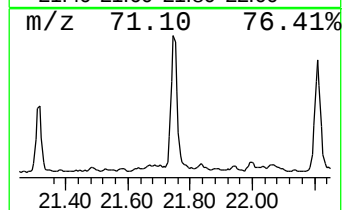
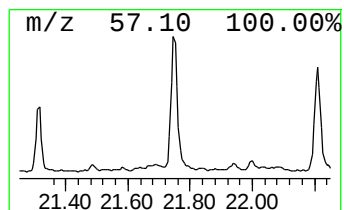
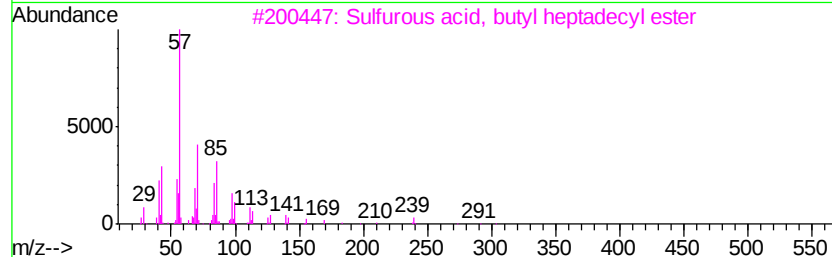
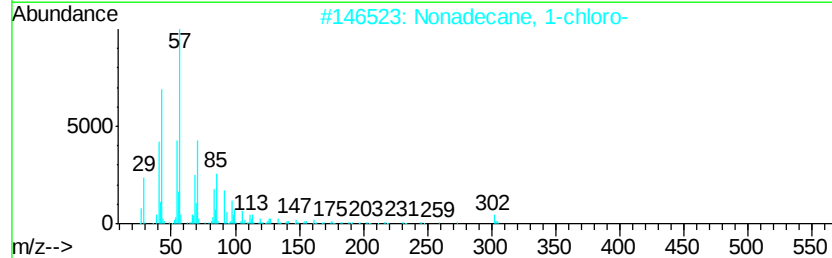
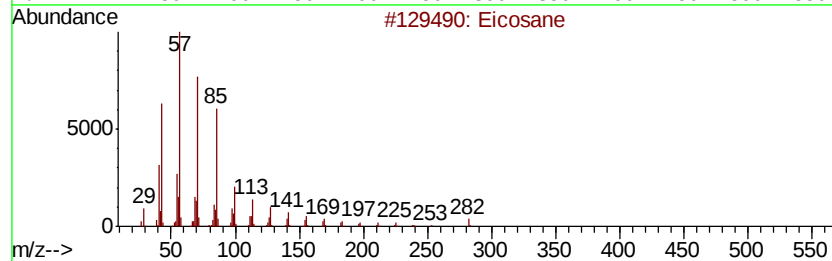
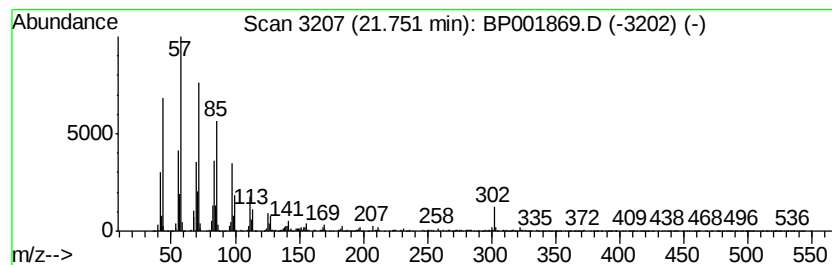
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.75	3.53 ng/ul	1130120	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	97
2		Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	94
3		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	91
4		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
5		Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
Operator : CG/JU
Sample : L1536-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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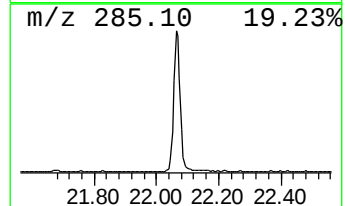
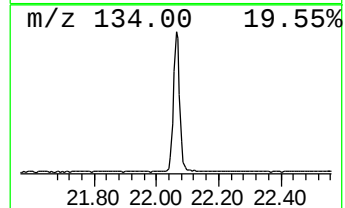
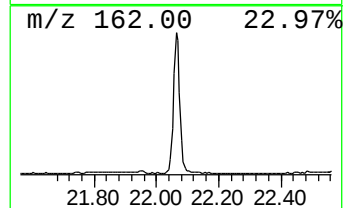
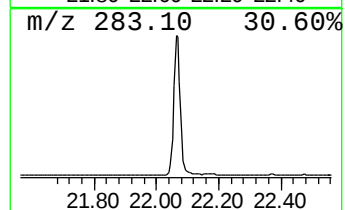
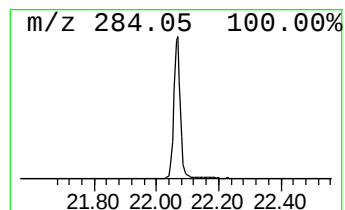
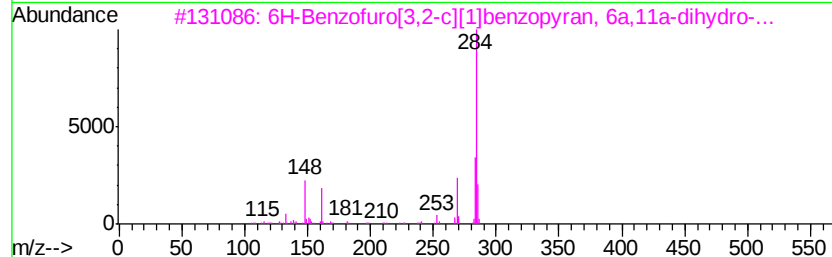
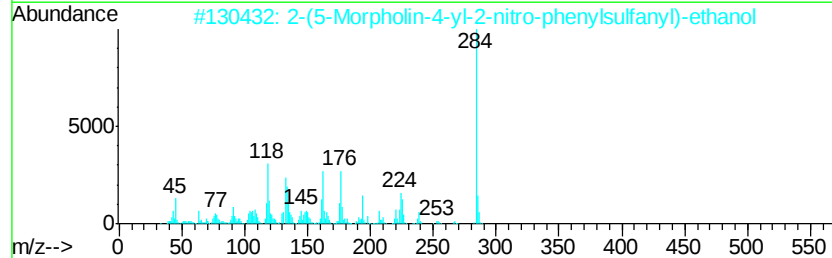
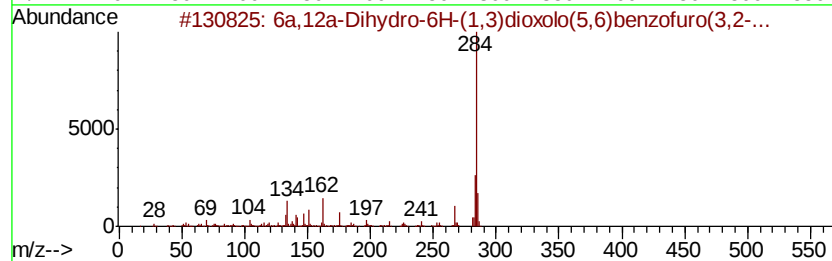
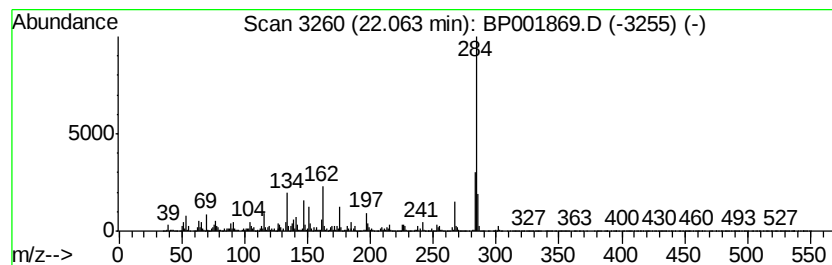
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 6a,12a-Dihydro-6H-(1,3)diox... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.06	12.30 ng/ul	3942340	Chrysene-d12	21.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6a,12a-Dihydro-6H-(1,3)dioxolo(5...	284	C16H12O5	022091-18-5	95	
2	2-(5-Morpholin-4-yl-2-nitro-phen...	284	C12H16N2O4S	1000294-41-6	53	
3	6H-Benzofuro[3,2-cl[1]benzopyran...	284	C17H16O4	000606-91-7	53	
4	4-Amino-5-(4-nitrophenylazo)benz...	284	C12H8N6O3	166766-07-0	50	
5	Benzofuran-3-one, 2-[3-hydroxy-4...	284	C16H12O5	1000128-62-8	49	



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
Operator : CG/JU
Sample : L1536-13
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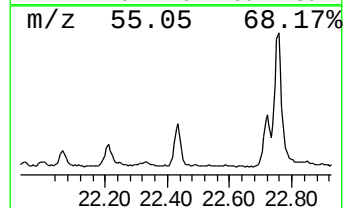
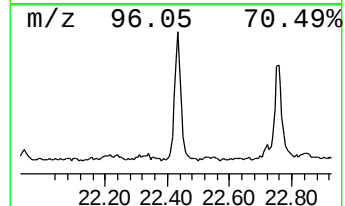
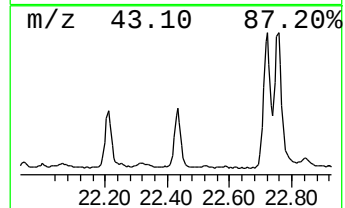
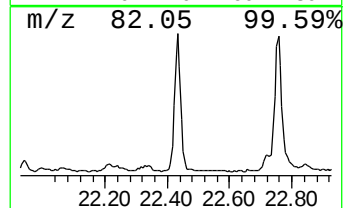
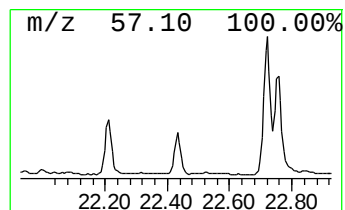
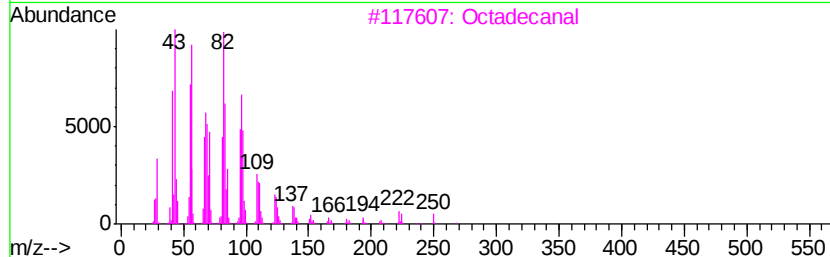
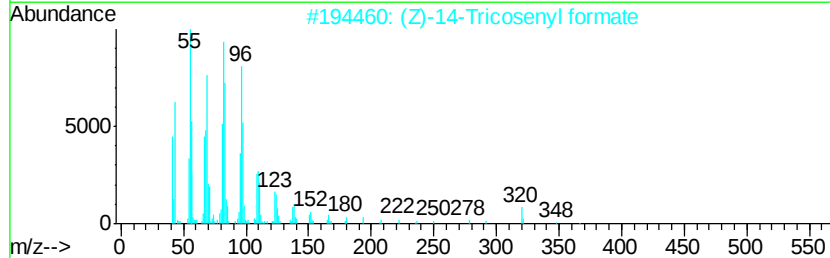
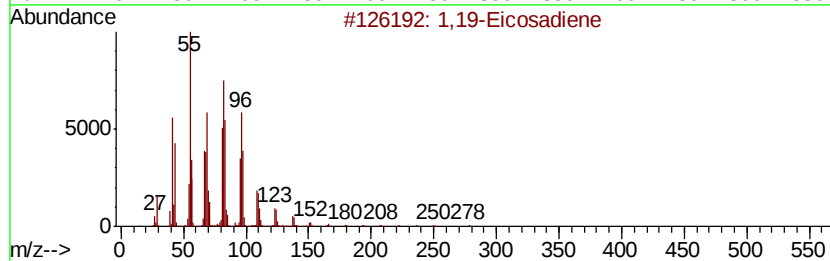
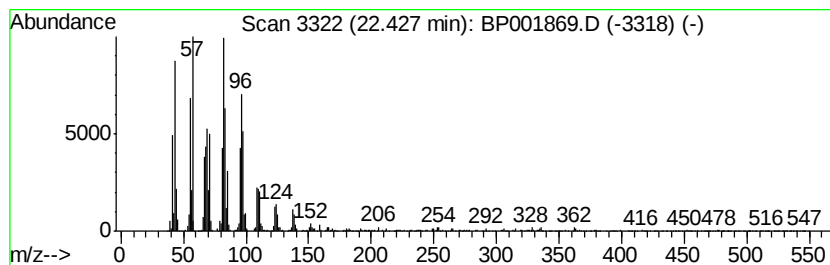
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,19-Eicosadiene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.43	4.85 ng/ul	1038010	Perylene-d12	23.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,19-Eicosadiene		278 C20H38	014811-95-1 94
2	(Z)-14-Tricosenyl formate		366 C24H46O2	077899-10-6 91
3	Octadecanal		268 C18H36O	000638-66-4 91
4	Octadecanal		268 C18H36O	000638-66-4 91
5	Oxirane, heptadecyl-		282 C19H38O	067860-04-2 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
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Sample : L1536-13
Misc :
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Instrument :
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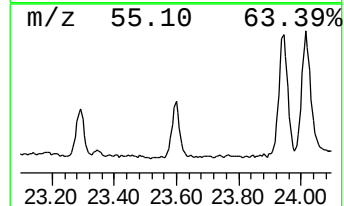
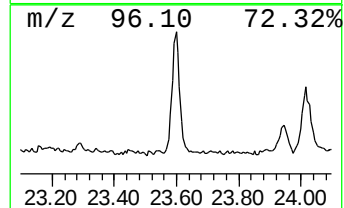
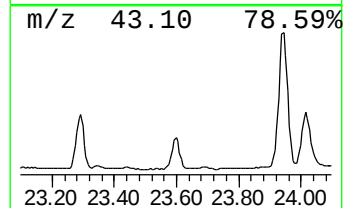
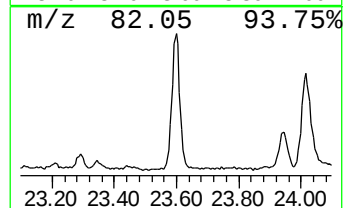
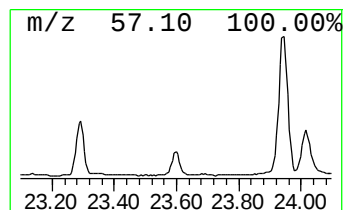
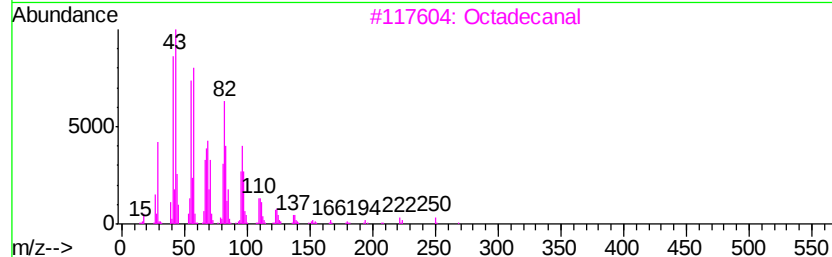
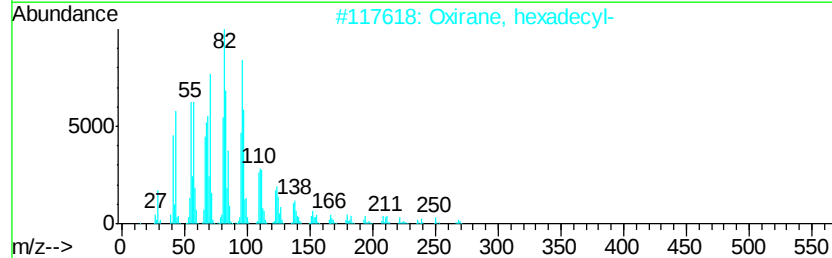
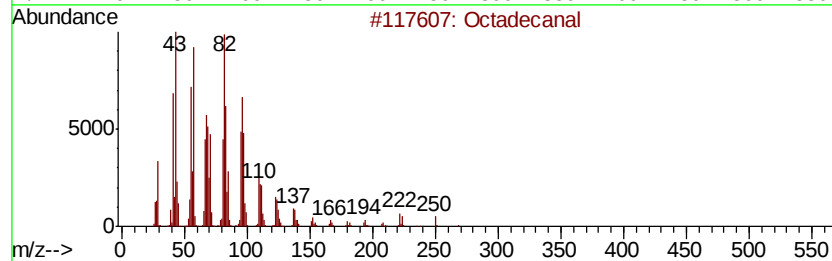
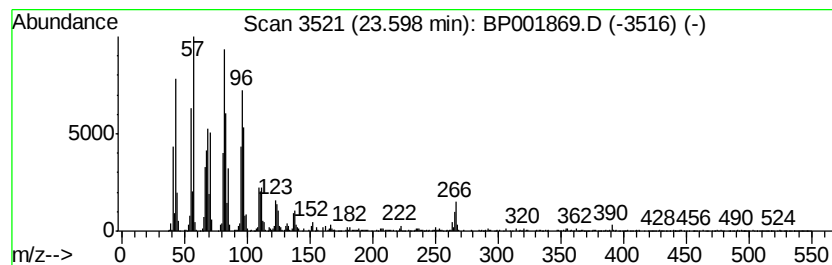
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.60	3.82 ng/ul	816846	Perylene-d12	23.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal		268	C18H36O	000638-66-4	93
2	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	93
3	Octadecanal		268	C18H36O	000638-66-4	90
4	Heptadecanal		254	C17H34O	1000376-70-0	87
5	Oxirane, hexadecyl-		268	C18H36O	007390-81-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
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Sample : L1536-13
Misc :
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ClientSampleId :
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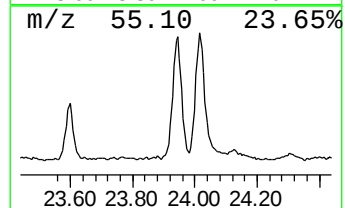
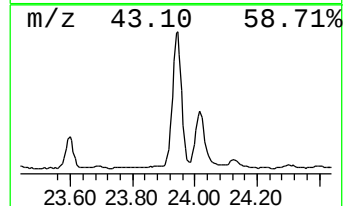
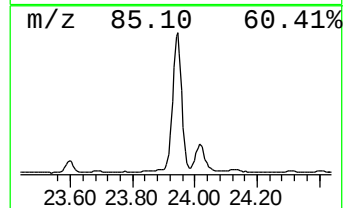
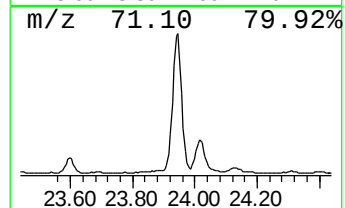
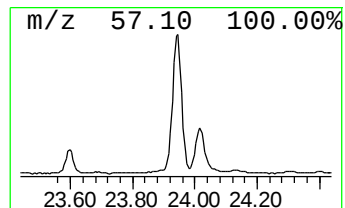
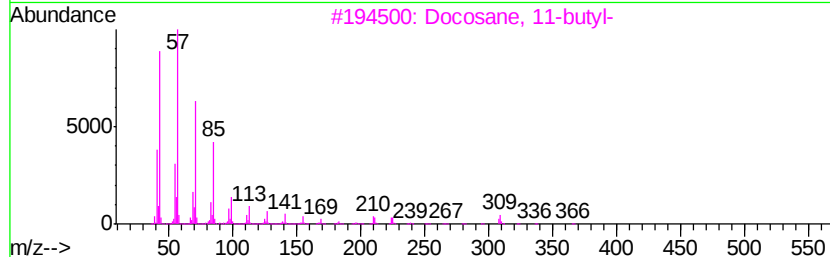
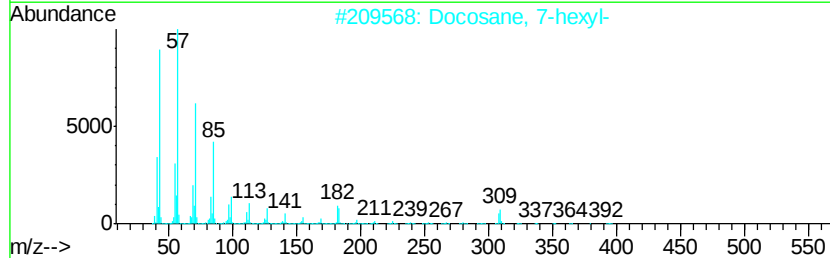
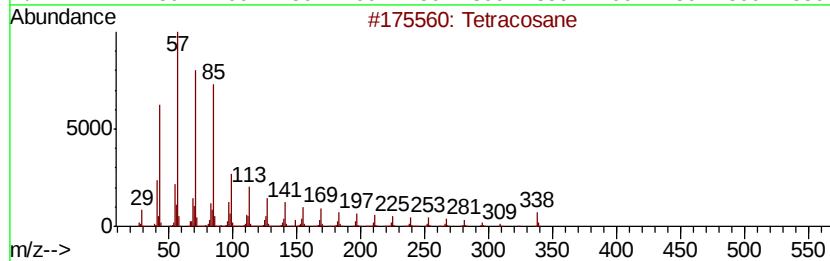
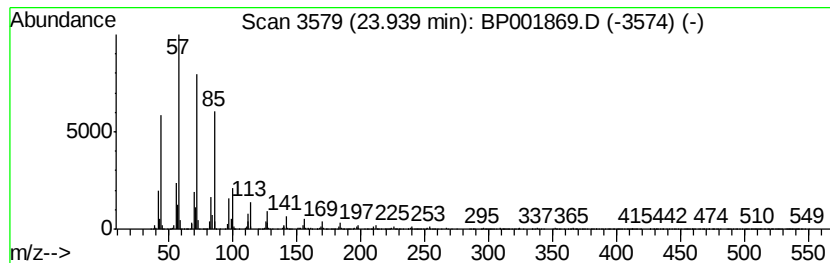
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.94	11.29 ng/ul	2415770	Perylene-d12	23.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	95
2		Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		triacontane	422	C30H62	000638-68-6	91



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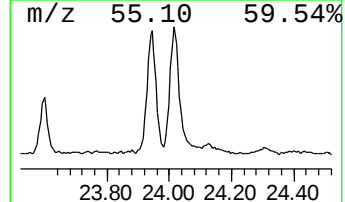
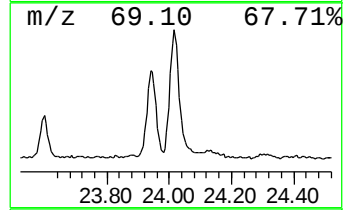
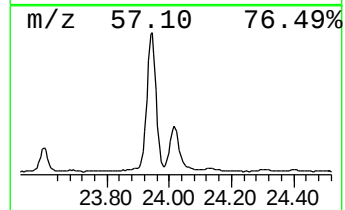
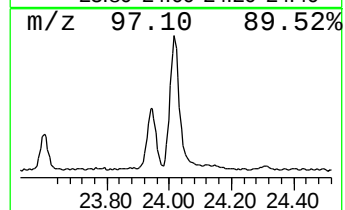
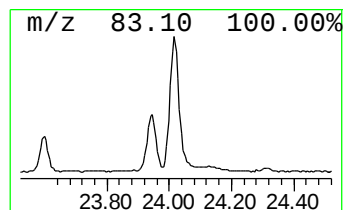
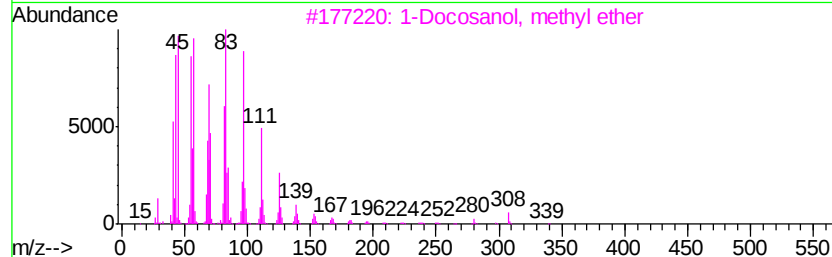
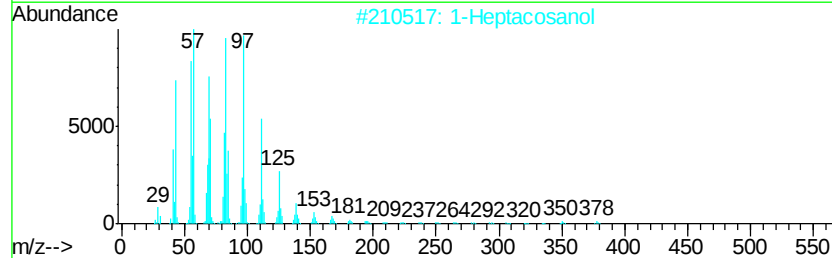
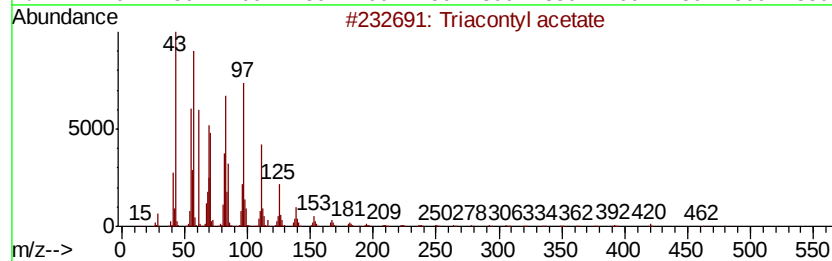
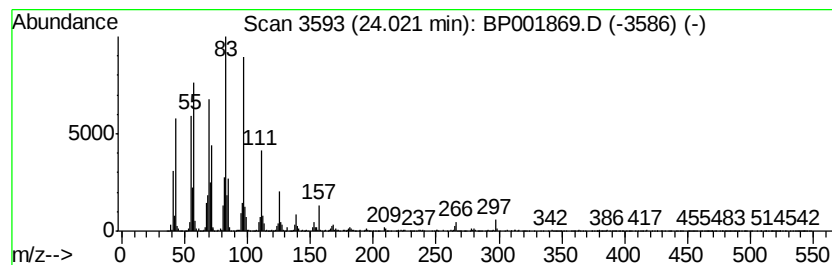
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 Quant Title : SVOA CALIBRATION

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

 Peak Number 9 Triacontyl acetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.02	7.91 ng/ul	1692820	Perylene-d12	23.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontyl acetate	480	C32H64O2	041755-58-2	96
2		1-Heptacosanol	396	C27H56O	002004-39-9	93
3		1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	93
4		Octacosanol	410	C28H58O	000557-61-9	81
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP022420\
Data File : BP001869.D
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ALS Vial : 17 Sample Multiplier: 1

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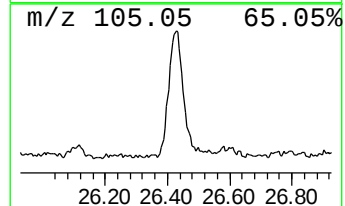
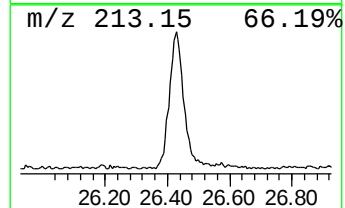
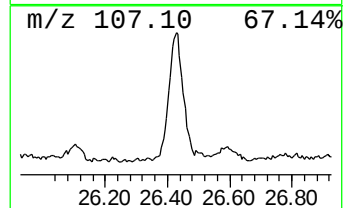
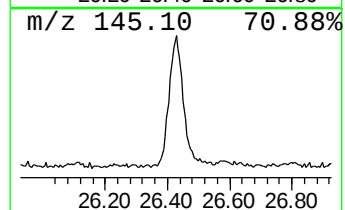
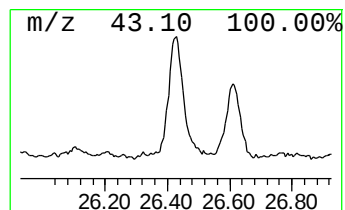
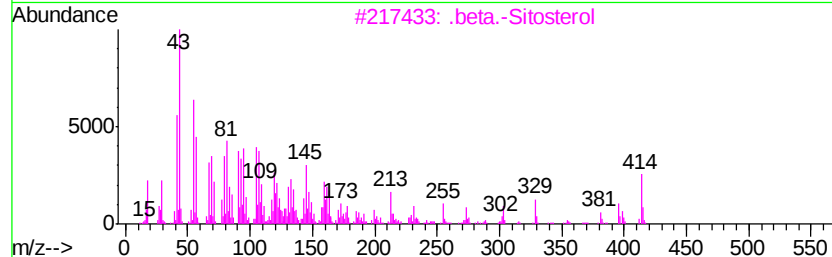
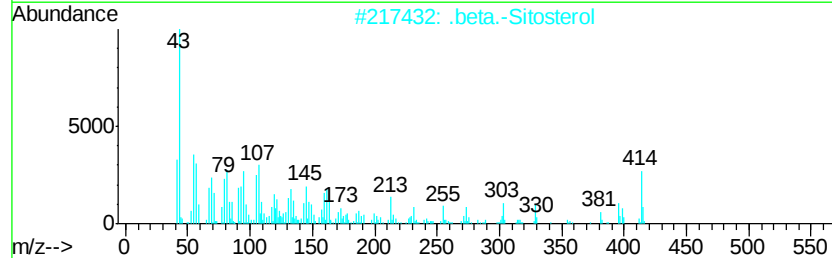
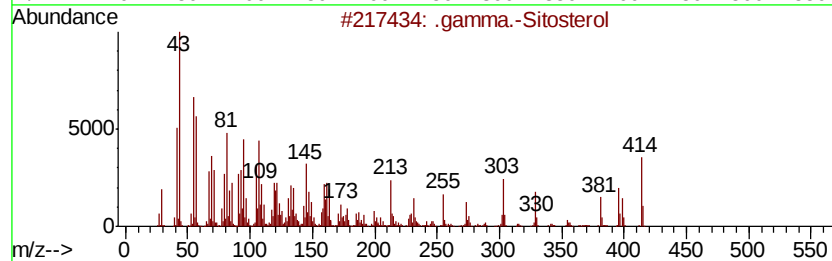
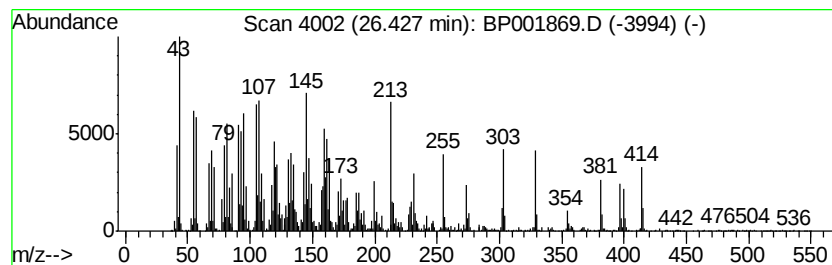
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 .gamma.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.43	10.89 ng/ul	2329170	Perylene-d12	23.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	89
3		.beta.-Sitosterol	414	C29H50O	000083-46-5	48
4		17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	25
5		cis-4-Acetoxy-trans-1-(m-methoxy...	273	C16H19NO3	1000241-10-4	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022420\
Data File : BP001869.D
Acq On : 24 Feb 2020 20:00
Operator : CG/JU
Sample : L1536-13
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB6X3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.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
Butane, 2-methoxy...	2.92	35.3	ng/ul	3476670	1	7.65	1969350	20.0
Nonanal	8.98	2.7	ng/ul	261890	1	7.65	1969350	20.0
1-Heneicosyl formate	20.87	5.6	ng/ul	1782030	5	21.13	6409190	20.0
(DEL) Alkane: Str...	21.75	3.5	ng/ul	1130120	5	21.13	6409190	20.0
6a,12a-Dihydro-6H...	22.06	12.3	ng/ul	3942340	5	21.13	6409190	20.0
1,19-Eicosadiene	22.43	4.8	ng/ul	1038010	6	23.35	4278810	20.0
Octadecanal	23.60	3.8	ng/ul	816846	6	23.35	4278810	20.0
(DEL) Alkane: Str...	23.94	11.3	ng/ul	2415770	6	23.35	4278810	20.0
Triacetyl acetate	24.02	7.9	ng/ul	1692820	6	23.35	4278810	20.0
.gamma.-Sitosterol	26.43	10.9	ng/ul	2329170	6	23.35	4278810	20.0