

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020309.D
 Acq On : 12 May 2019 21:30
 Operator : JU/SJ
 Sample : K2766-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS008-0002-01

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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_M\METHODS\8270-BM043019.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.981	164	169	178	rVB	216778	307569	3.14%	0.360%
2	5.346	395	401	409	rBV	1077089	1621864	16.57%	1.900%
3	6.934	663	671	686	rBV	1071258	1778778	18.17%	2.084%
4	7.293	725	732	747	rVB	1130708	1855896	18.96%	2.174%
5	7.763	806	812	821	rBV	290370	463317	4.73%	0.543%
6	8.081	859	866	877	rBV	799937	1278911	13.07%	1.498%
7	8.928	1002	1010	1020	rBV	657726	1139800	11.65%	1.335%
8	10.551	1280	1286	1298	rVB	326856	573356	5.86%	0.672%
9	13.022	1698	1706	1718	rBV	1404338	2165721	22.13%	2.537%
10	13.339	1756	1760	1770	rVB3	58765	107678	1.10%	0.126%
11	13.869	1844	1850	1862	rVB2	270475	519483	5.31%	0.609%
12	14.410	1936	1942	1950	rBV2	539972	833141	8.51%	0.976%
13	15.904	2190	2196	2201	rBV	1297040	1877134	19.18%	2.199%
14	16.327	2264	2268	2273	rBV	133252	149650	1.53%	0.175%
15	16.551	2301	2306	2314	rBV2	216540	357933	3.66%	0.419%
16	17.074	2392	2395	2404	rVV3	56581	109164	1.12%	0.128%
17	17.157	2404	2409	2418	rVV	713118	1263340	12.91%	1.480%
18	17.233	2419	2422	2429	rVB2	124668	160345	1.64%	0.188%
19	17.321	2432	2437	2441	rBV2	114180	159581	1.63%	0.187%
20	17.527	2470	2472	2481	rVB6	63803	98590	1.01%	0.116%
21	17.633	2484	2490	2495	rBV2	147654	215914	2.21%	0.253%
22	17.774	2509	2514	2519	rBV	169595	244201	2.50%	0.286%
23	17.921	2534	2539	2547	rBV3	86526	163101	1.67%	0.191%
24	18.045	2555	2560	2571	rBV	882711	1343073	13.72%	1.574%
25	18.227	2586	2591	2597	rVB3	101352	141993	1.45%	0.166%
26	18.286	2597	2601	2604	rBV	80576	99769	1.02%	0.117%
27	18.321	2604	2607	2615	rVV5	113668	232814	2.38%	0.273%
28	18.509	2635	2639	2644	rBV5	80548	111183	1.14%	0.130%
29	18.562	2644	2648	2653	rVV2	97710	128668	1.31%	0.151%
30	19.051	2727	2731	2733	rBV	191362	263365	2.69%	0.309%
31	19.074	2733	2735	2744	rVB3	237745	303012	3.10%	0.355%
32	19.215	2755	2759	2763	rBV2	96473	178117	1.82%	0.209%
33	19.262	2763	2767	2773	rVV2	215555	296110	3.03%	0.347%
34	19.327	2774	2778	2780	rVV	292326	345589	3.53%	0.405%

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35	19.356	2780	2783	2790	rVB2	367573	484820	4.95%	0.568%
36	19.586	2818	2822	2824	rBV	98308	127668	1.30%	0.150%
37	19.786	2851	2856	2863	rBV	2697763	3555448	36.33%	4.166%
38	19.927	2876	2880	2884	rVB	280871	314193	3.21%	0.368%
39	19.968	2884	2887	2894	rVB2	127153	207645	2.12%	0.243%
40	20.062	2898	2903	2910	rBV3	1009766	1895524	19.37%	2.221%
41	20.262	2934	2937	2944	rVB2	160114	215733	2.20%	0.253%
42	20.345	2946	2951	2955	rBV	1050261	1291595	13.20%	1.513%
43	20.403	2955	2961	2972	rVV6	269091	597095	6.10%	0.700%
44	20.503	2973	2978	2984	rVV3	378706	594689	6.08%	0.697%
45	20.580	2986	2991	2997	rVB5	159946	330703	3.38%	0.387%
46	20.662	2997	3005	3010	rBV	617948	1148125	11.73%	1.345%
47	20.756	3018	3021	3025	rVV2	106958	123665	1.26%	0.145%
48	20.803	3025	3029	3034	rVB	535782	641162	6.55%	0.751%
49	20.868	3037	3040	3044	rBV	222513	247098	2.52%	0.290%
50	21.045	3065	3070	3073	rBV	4322745	4796123	49.00%	5.619%
51	21.074	3073	3075	3080	rVV2	688816	832819	8.51%	0.976%
52	21.133	3080	3085	3090	rVB2	373716	561562	5.74%	0.658%
53	21.192	3092	3095	3097	rBV3	93183	101497	1.04%	0.119%
54	21.221	3097	3100	3104	rVV3	93144	102076	1.04%	0.120%
55	21.350	3116	3122	3131	rBV2	1453314	3085500	31.53%	3.615%
56	21.486	3141	3145	3147	rBV2	314298	436542	4.46%	0.511%
57	21.674	3172	3177	3183	rBV2	484255	796670	8.14%	0.933%
58	21.733	3184	3187	3192	rVB2	191467	252519	2.58%	0.296%
59	21.803	3195	3199	3205	rBV3	231223	325076	3.32%	0.381%
60	21.945	3214	3223	3232	rBV2	5106923	9787188	100.00%	11.467%
61	22.015	3233	3235	3241	rVB	192570	247323	2.53%	0.290%
62	22.386	3293	3298	3302	rBV2	576945	887509	9.07%	1.040%
63	22.433	3302	3306	3312	rVV	414340	656030	6.70%	0.769%
64	22.521	3316	3321	3325	rVB	564641	797229	8.15%	0.934%
65	22.627	3335	3339	3347	rVB	355078	504127	5.15%	0.591%
66	22.909	3381	3387	3391	rBV	1314377	2104124	21.50%	2.465%
67	22.962	3391	3396	3405	rVV	1969906	3170781	32.40%	3.715%
68	23.044	3406	3410	3417	rVB	359686	528475	5.40%	0.619%
69	23.486	3480	3485	3490	rVB	322266	551251	5.63%	0.646%
70	23.639	3505	3511	3519	rVB	693177	1477036	15.09%	1.731%
71	23.809	3536	3540	3548	rVB	407267	769728	7.86%	0.902%

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72	24.050	3577	3581	3586	rBV2	223973	430734	4.40%	0.505%
73	24.144	3592	3597	3607	rVB	727326	1491358	15.24%	1.747%
74	24.244	3607	3614	3625	rVV	908567	1904044	19.45%	2.231%
75	24.350	3627	3632	3643	rVB	456617	968976	9.90%	1.135%
76	24.521	3656	3661	3673	rVB5	426506	1016934	10.39%	1.191%
77	24.639	3676	3681	3688	rVV	255653	533064	5.45%	0.625%
78	25.380	3803	3807	3818	rVB3	331078	762542	7.79%	0.893%
79	25.968	3903	3907	3920	rVB6	292989	979665	10.01%	1.148%
80	26.109	3923	3931	3942	rVB4	629333	1978963	20.22%	2.319%
81	26.750	4033	4040	4053	rVB5	478332	1551966	15.86%	1.818%
82	26.944	4066	4073	4085	rVB5	737347	2308380	23.59%	2.705%
83	27.338	4133	4140	4157	rVB6	548607	2257631	23.07%	2.645%
84	27.609	4179	4186	4194	rBV3	540606	1524410	15.58%	1.786%
85	28.079	4260	4266	4281	rVB5	657653	2209014	22.57%	2.588%

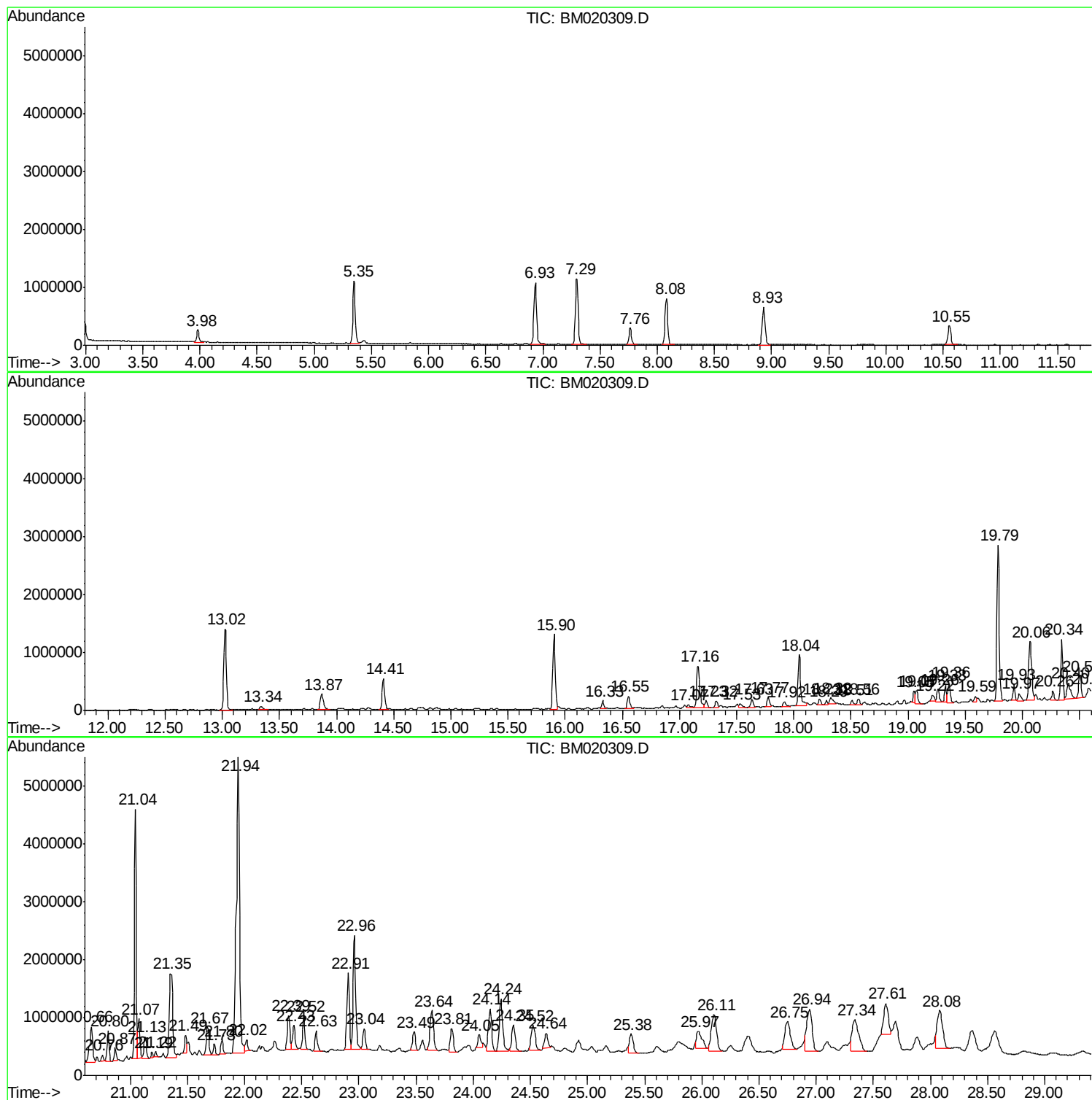
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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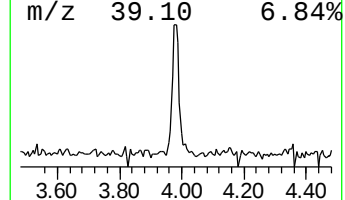
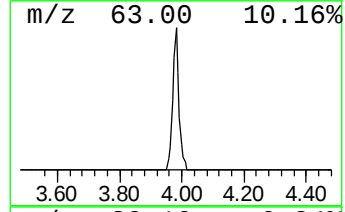
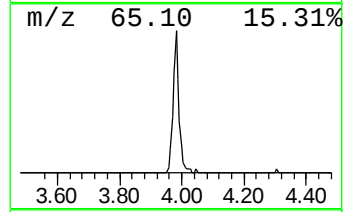
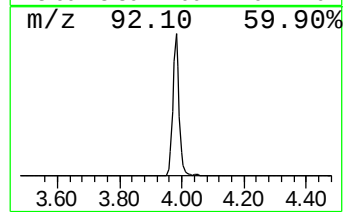
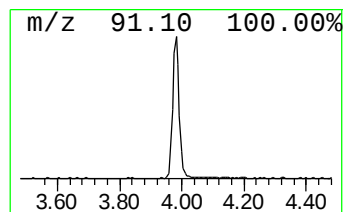
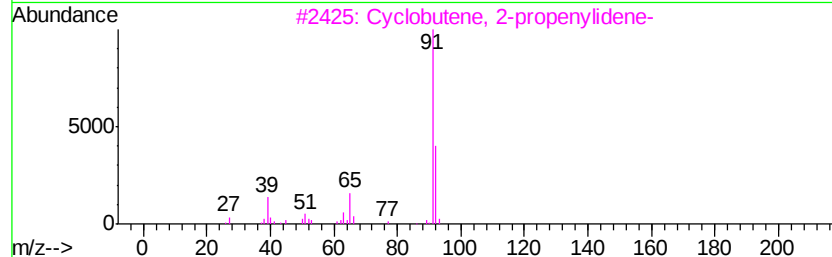
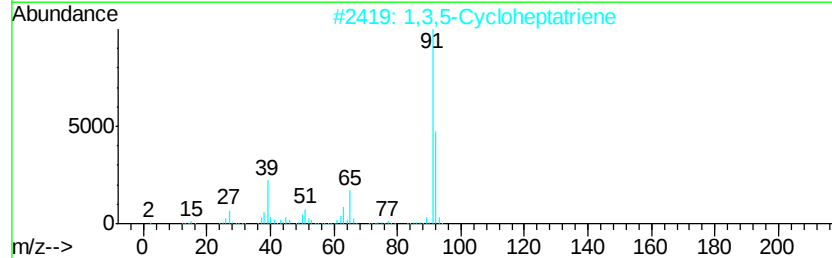
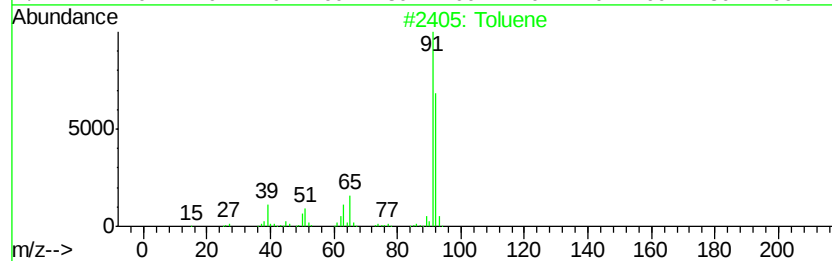
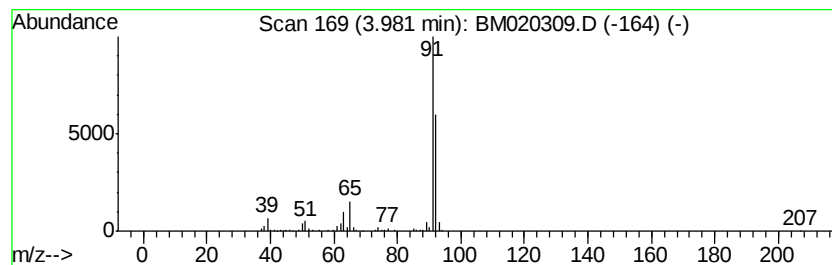
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Toluene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.98	13.28 ng	307569	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Toluene	92	C7H8	000108-88-3	91
2		1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	87
3		Cyclobutene, 2-propenylidene-	92	C7H8	052097-85-5	78
4		Molybdenum, di-.mu.-chlorobis[(1...	532	C20H26Cl2Mo2	035625-66-2	72
5		2,5-Norbornadiene	92	C7H8	000121-46-0	64



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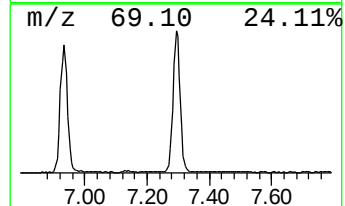
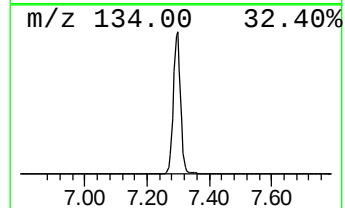
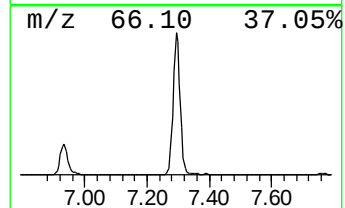
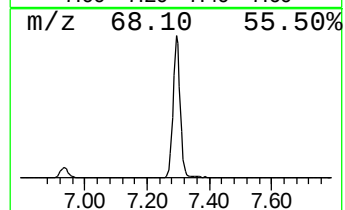
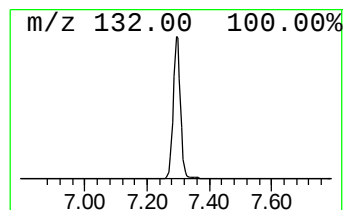
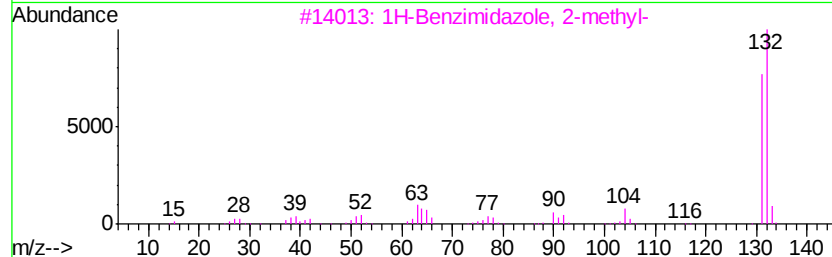
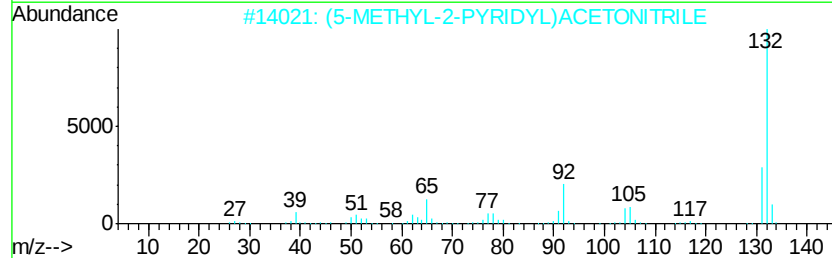
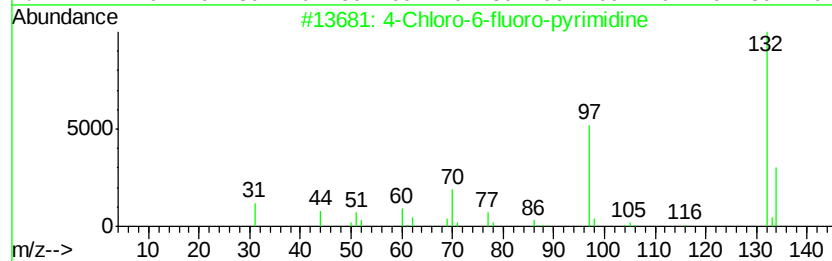
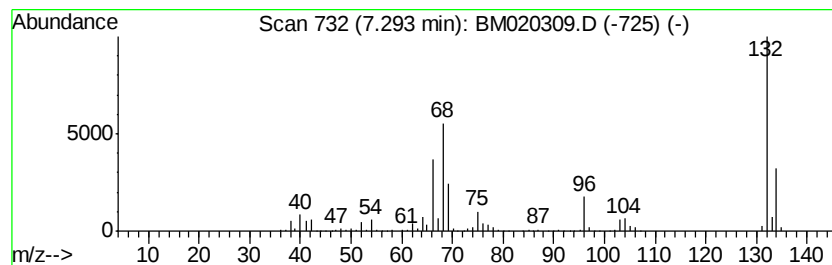
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.29 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	80.11 ng	1855900	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	22
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	13



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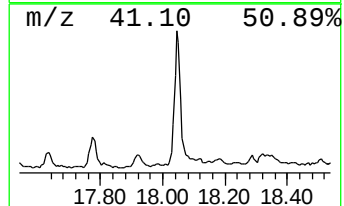
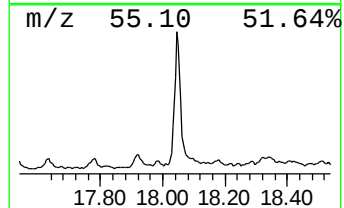
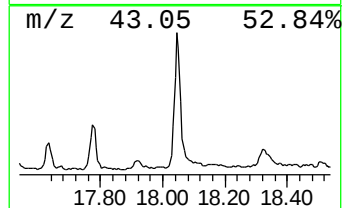
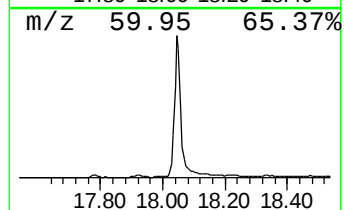
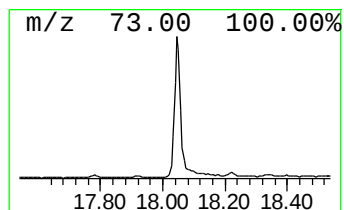
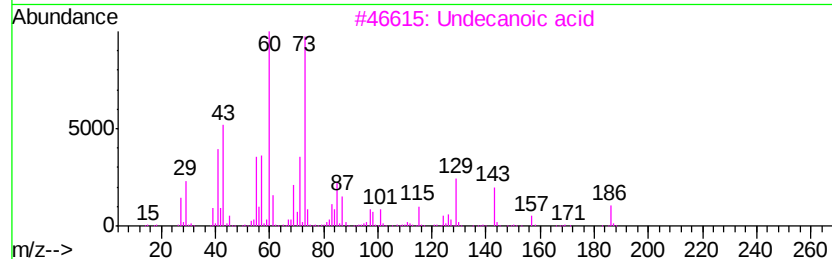
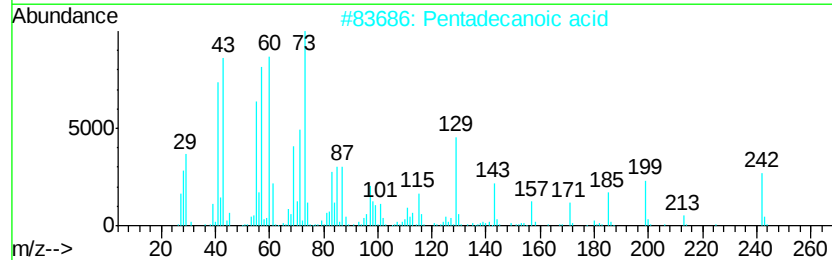
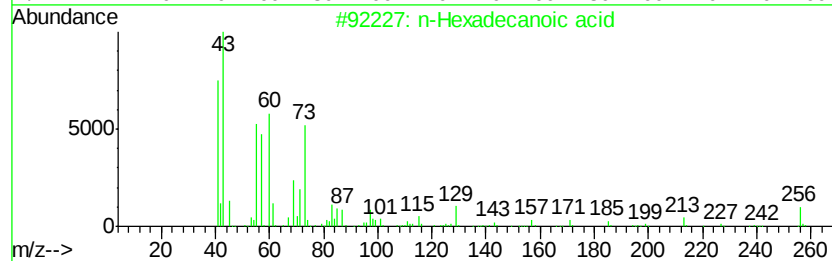
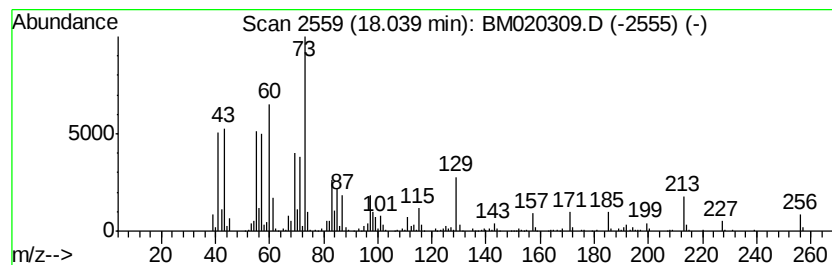
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	21.26 ng	1343070	Phenanthrene-d10	17.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
3		Undecanoic acid	186	C11H22O2	000112-37-8	46
4		Tridecanoic acid	214	C13H26O2	000638-53-9	45
5		Decane, 1-iodo-	268	C10H21I	002050-77-3	18



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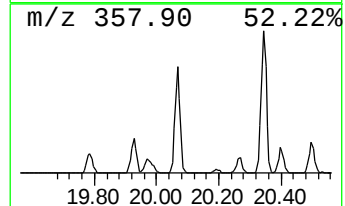
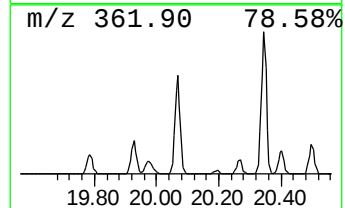
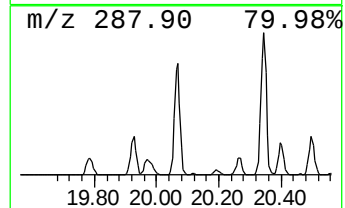
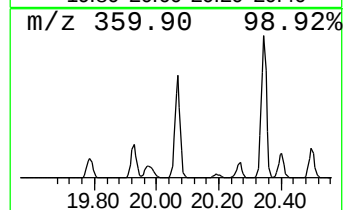
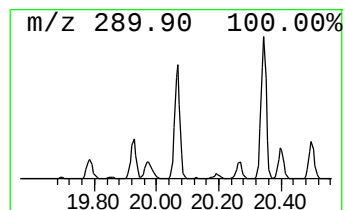
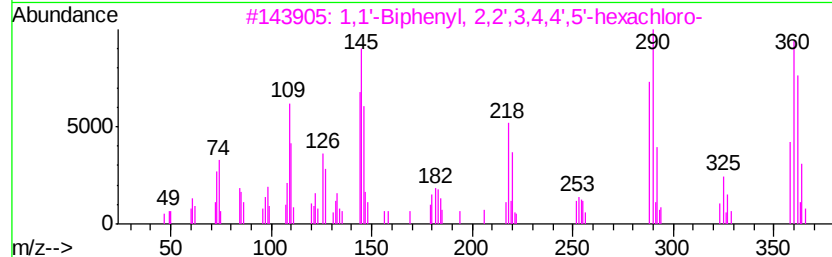
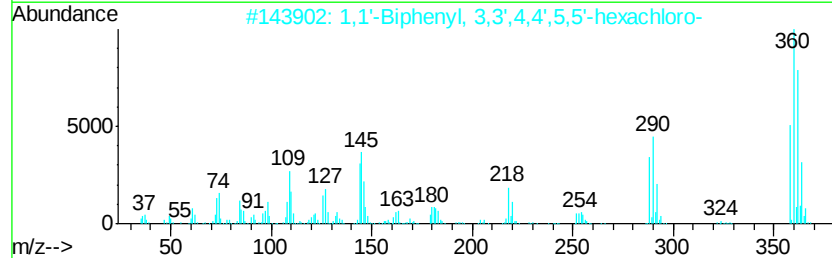
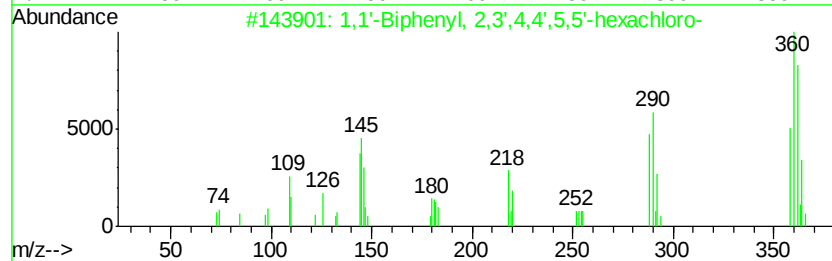
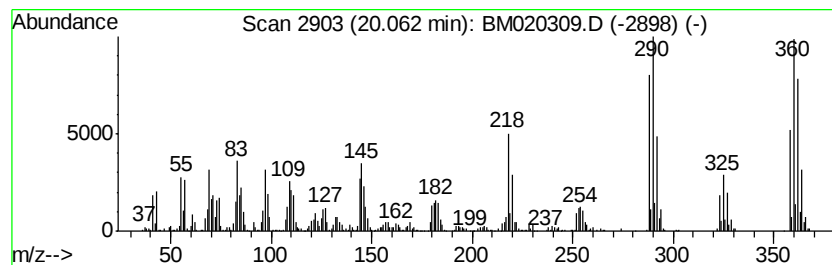
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1,1'-Biphenyl, 2,3',4,4',5,... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.06	12.29 ng	1895520	Chrysene-d12	21.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	99
2		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
3		1,1'-Biphenyl, 2,2',3,4,4',5'-he...	358	C12H4Cl6	035065-28-2	99
4		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99
5		1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

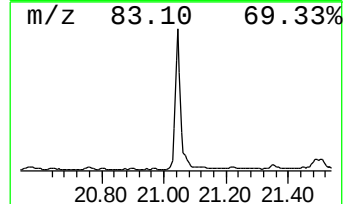
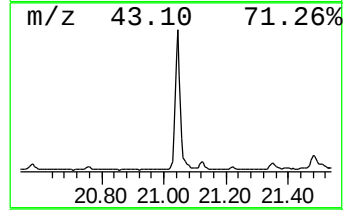
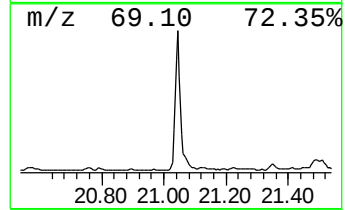
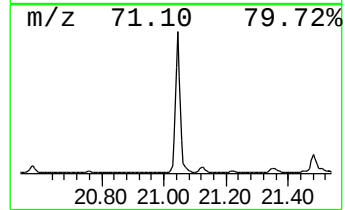
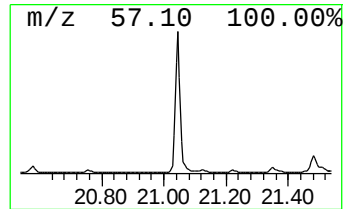
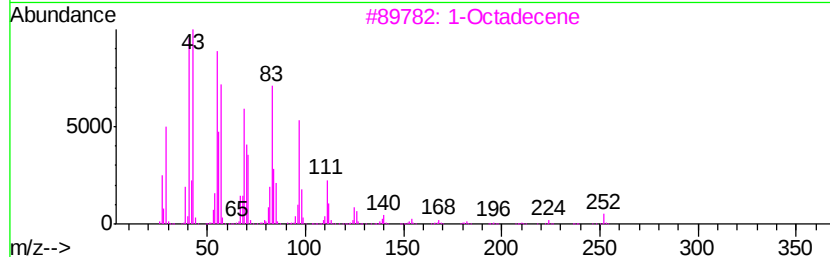
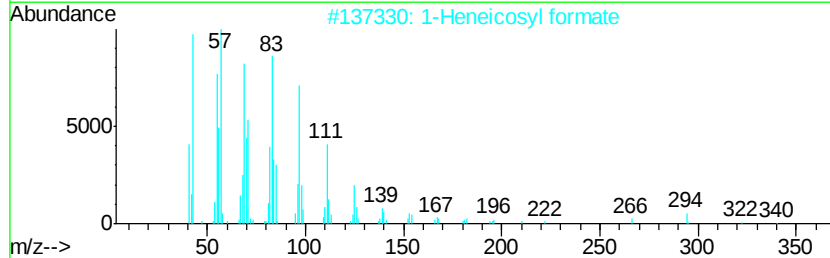
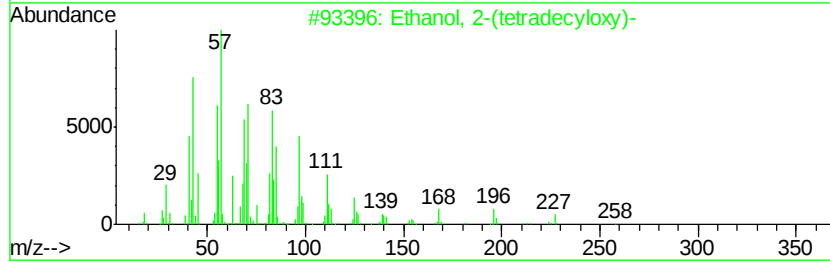
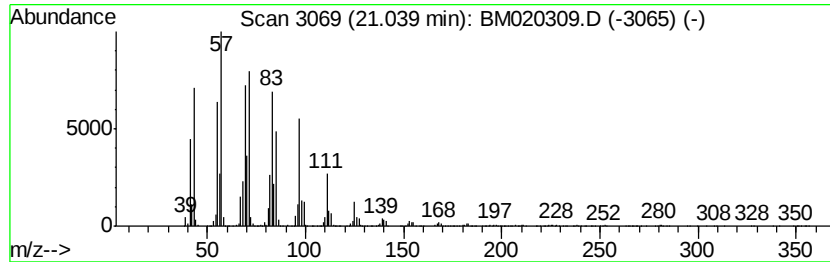
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.04	31.09 ng	4796120	Chrysene-d12	21.35		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	96
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	92
3		1-Octadecene	252	C18H36	000112-88-9	89
4		Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	83
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
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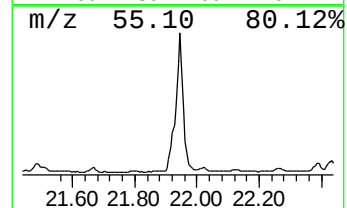
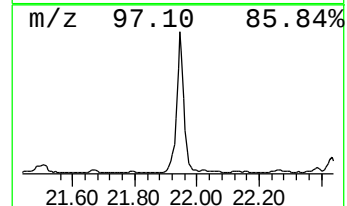
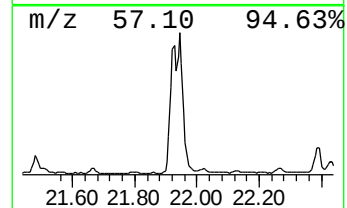
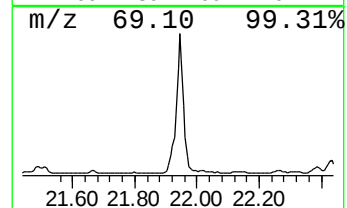
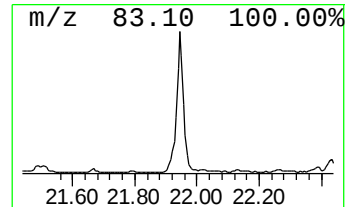
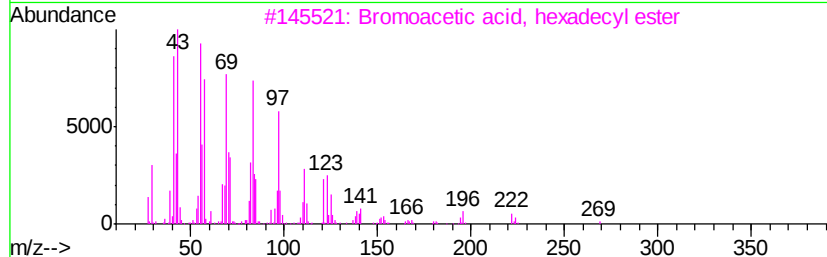
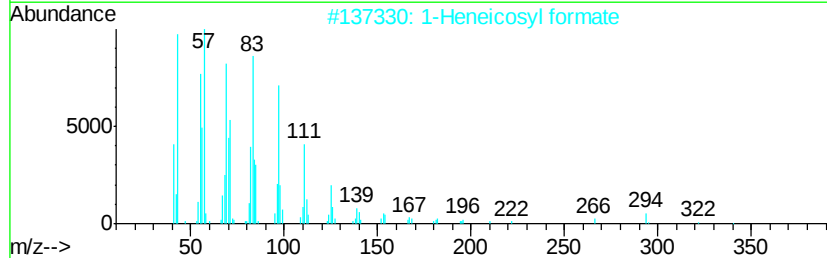
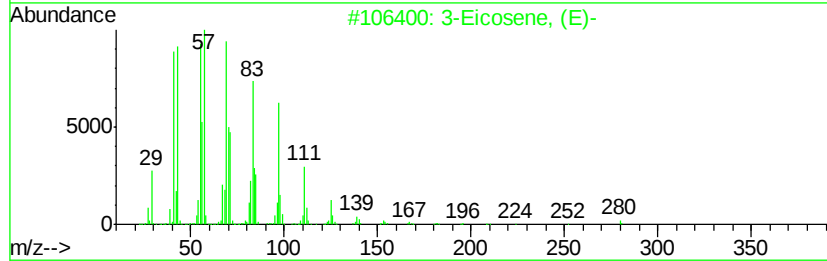
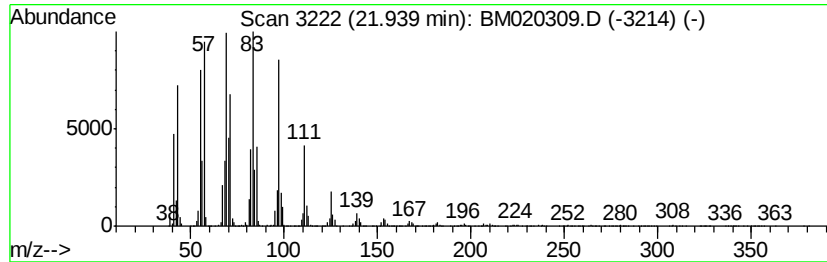
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 3-Eicosene, (E)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.94	63.44 ng	9787190	Chrysene-d12	21.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		3-Eicosene, (E)-	280 C20H40	074685-33-9 91
2		1-Heneicosyl formate	340 C22H44O2	077899-03-7 91
3		Bromoacetic acid, hexadecyl ester	362 C18H35BrO2	005454-48-8 86
4		1-Heptadecanol	256 C17H36O	001454-85-9 80
5		Dichloroacetic acid, tridecyl ester	310 C15H28Cl2O2	1000280-48-3 80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
 Data File : BM020309.D
 Acq On : 12 May 2019 21:30
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Instrument :
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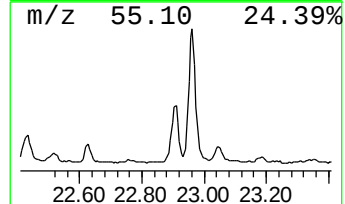
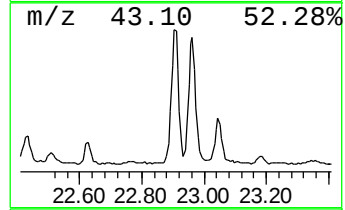
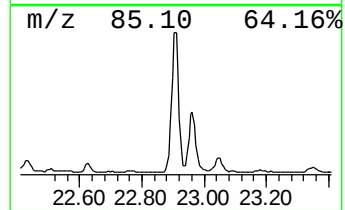
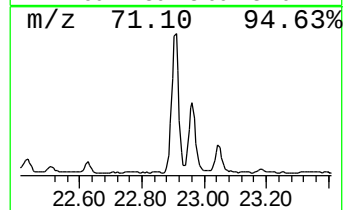
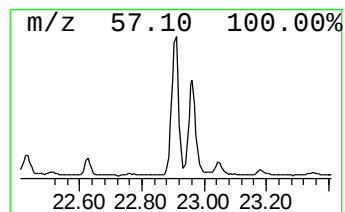
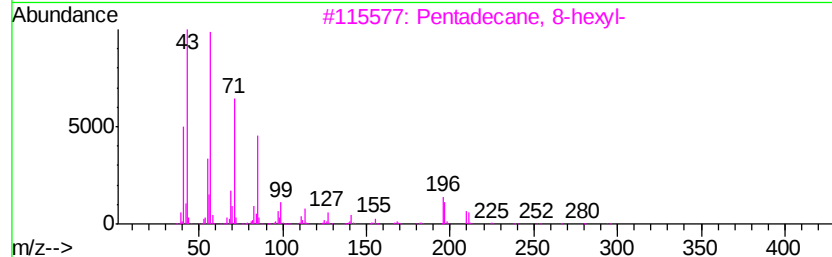
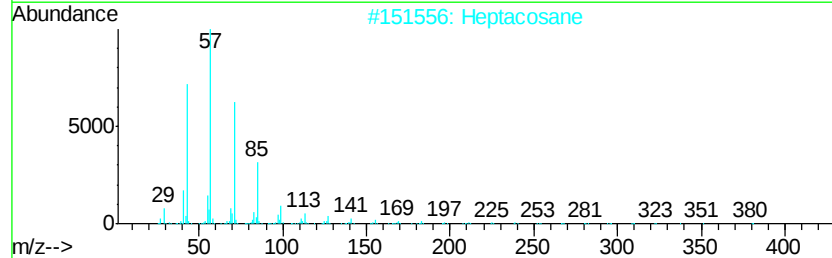
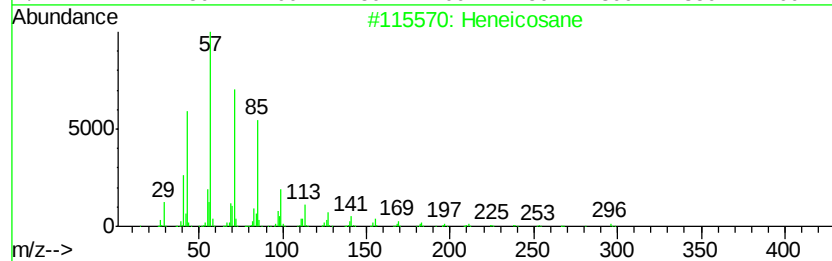
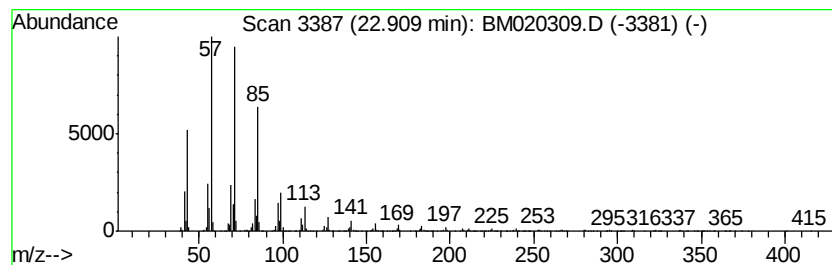
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 8 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.91	28.49 ng	2104120	Perylene-d12	23.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	90
4	Nonacosane		408	C29H60	000630-03-5	86
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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Sample : K2766-10
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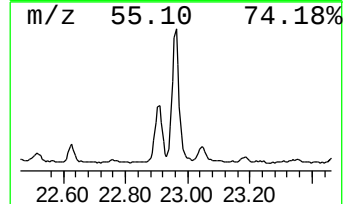
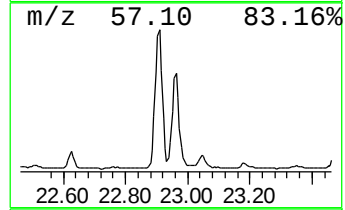
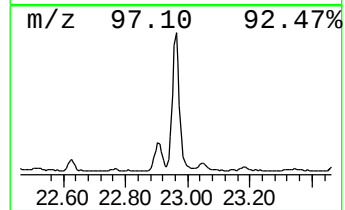
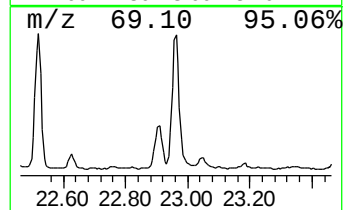
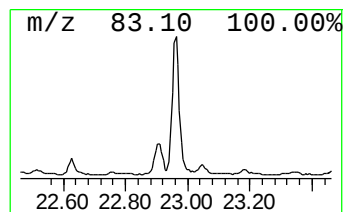
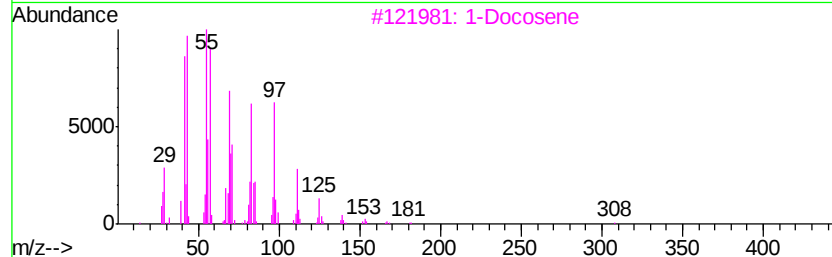
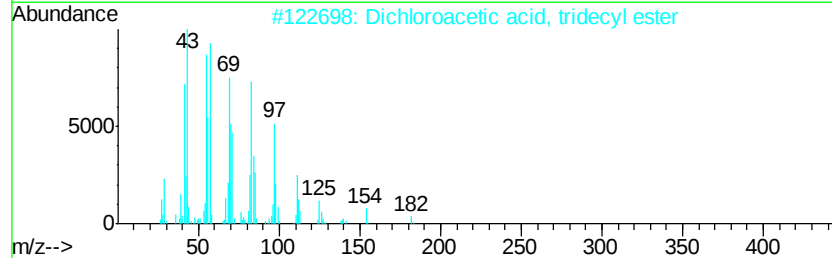
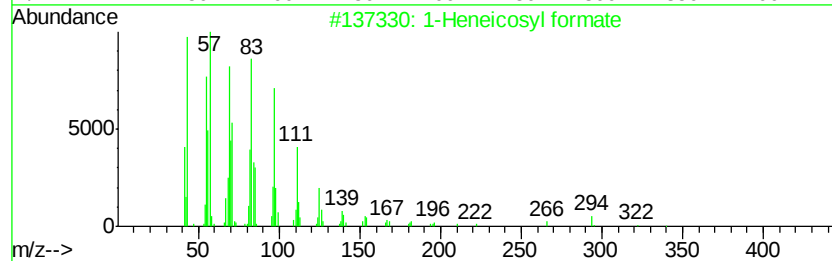
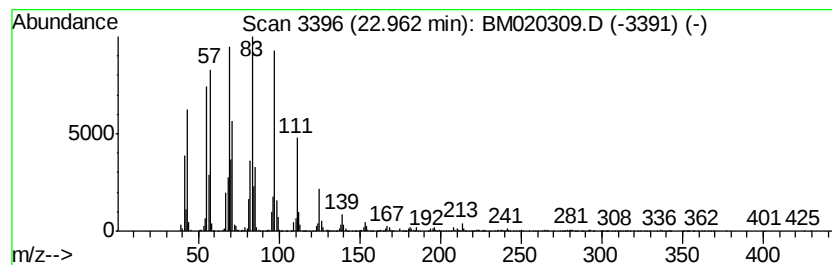
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.96	42.93 ng	3170780	Perylene-d12	23.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	87
2	Dichloroacetic acid, tridecyl ester	310	C15H28Cl2O2	1000280-48-3	86
3	1-Docosene	308	C22H44	001599-67-3	76
4	1-Tetracosanol	354	C24H50O	000506-51-4	64
5	1-Hexacosanol	382	C26H54O	000506-52-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
Misc :
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Instrument :
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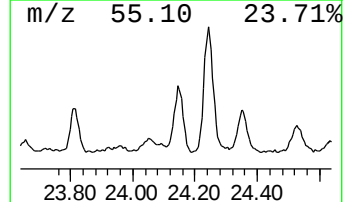
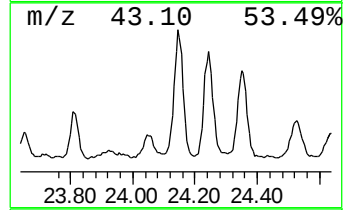
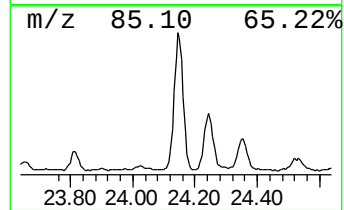
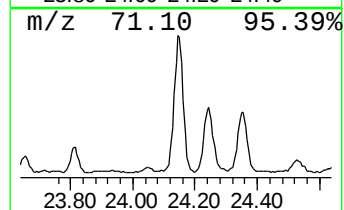
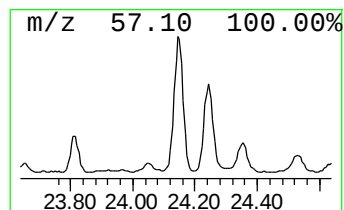
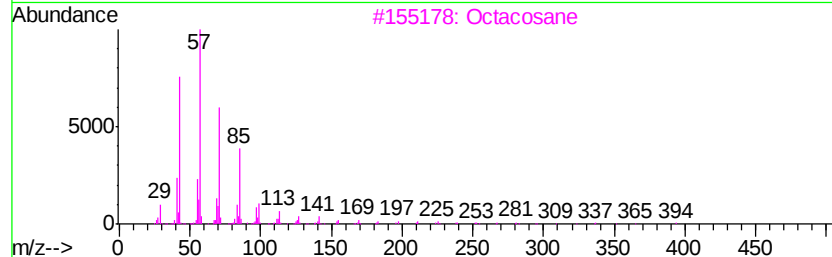
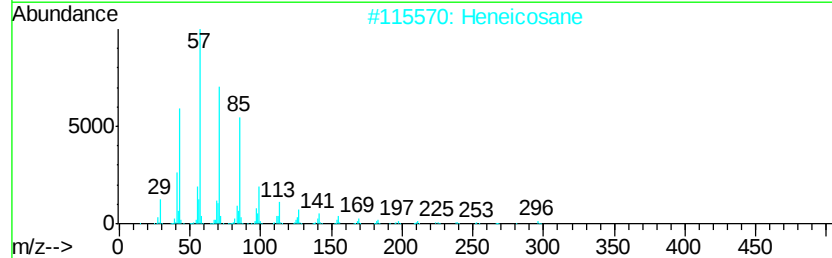
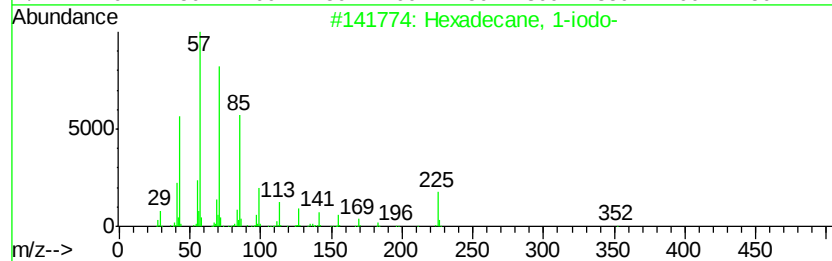
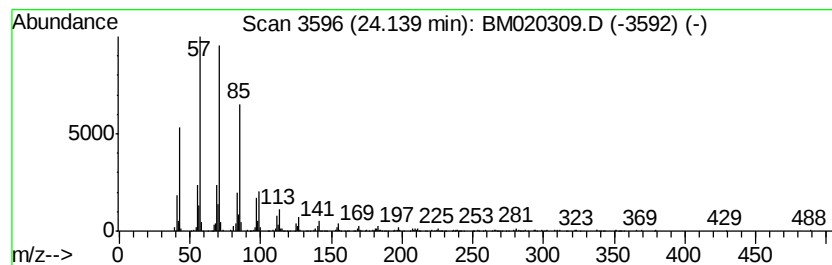
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Hexadecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.14	20.19 ng	1491360	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	92
2		Heneicosane	296	C21H44	000629-94-7	87
3		Octacosane	394	C28H58	000630-02-4	83
4		Heptacosane	380	C27H56	000593-49-7	74
5		Tetratetracontane	619	C44H90	007098-22-8	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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Sample : K2766-10
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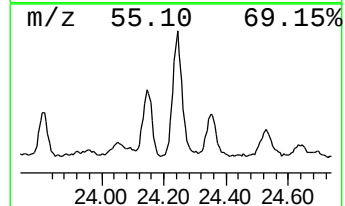
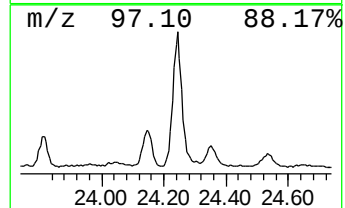
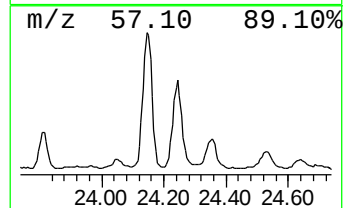
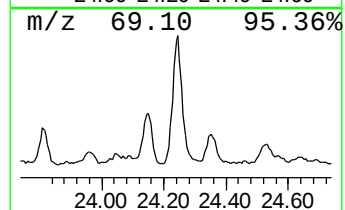
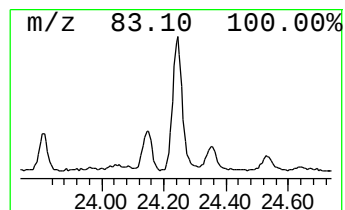
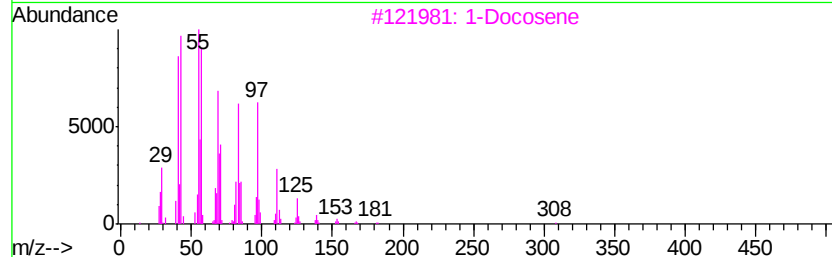
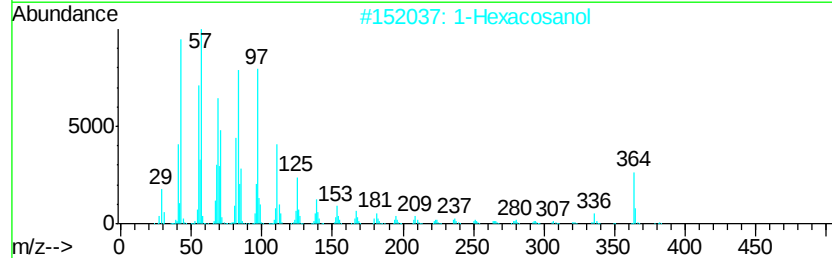
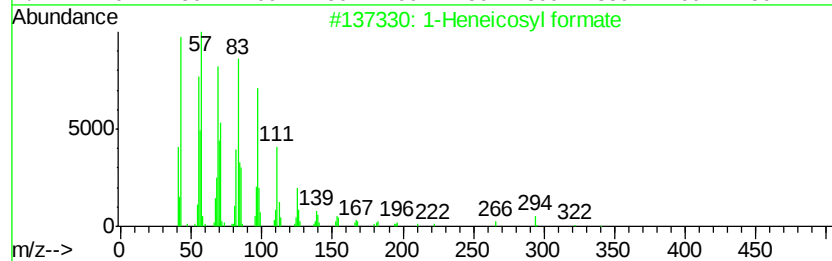
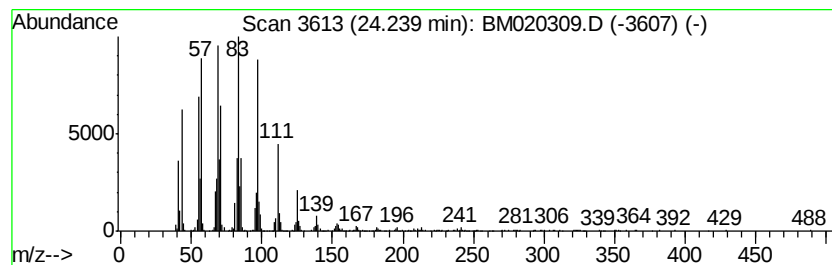
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Hexacosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.24	25.78 ng	1904040	Perylene-d12	23.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2	1-Hexacosanol	382	C26H54O	000506-52-5	76
3	1-Docosene	308	C22H44	001599-67-3	62
4	1-Tetracosanol	354	C24H50O	000506-51-4	58
5	17-Pentatriacontene	491	C35H70	006971-40-0	58



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Data File : BM020309.D
Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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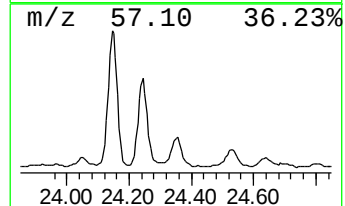
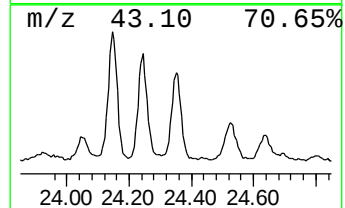
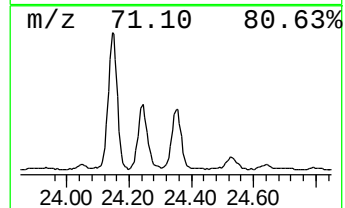
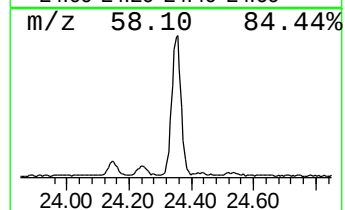
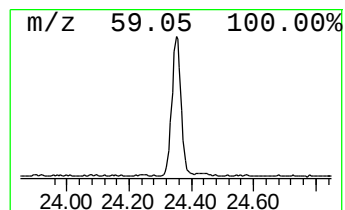
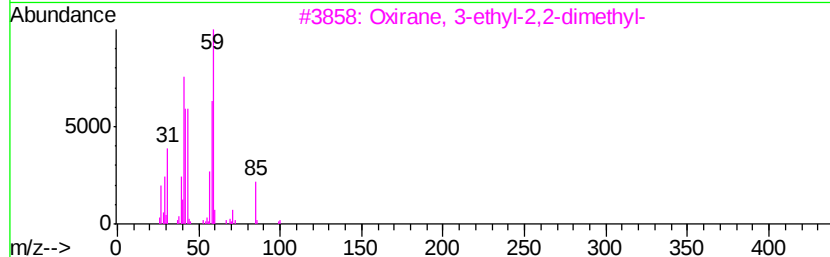
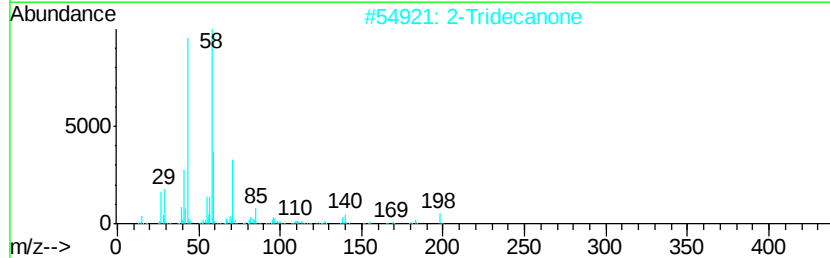
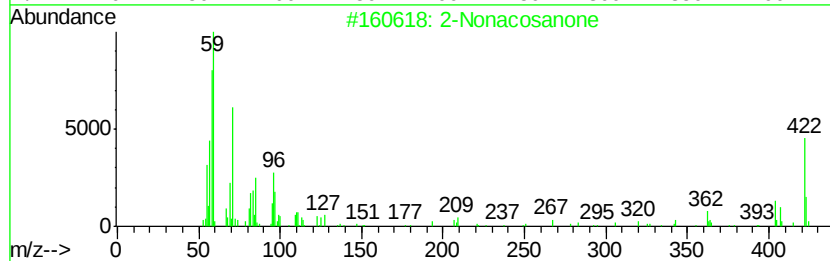
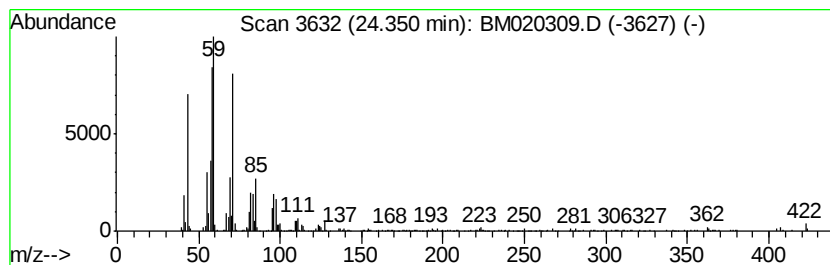
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 2-Nonacosanone Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.35	13.12 ng	968976	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Nonacosanone	422	C29H58O	017600-99-6	91
2		2-Tridecanone	198	C13H26O	000593-08-8	47
3		Oxirane, 3-ethyl-2,2-dimethyl-	100	C6H12O	001192-22-9	47
4		d-Ribo-tetrofuranose, 4-c-cyclop...	200	C10H16O4	030571-50-7	43
5		2-Nonadecanone	282	C19H38O	000629-66-3	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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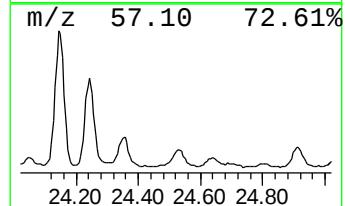
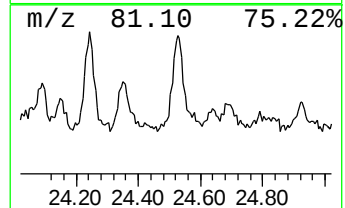
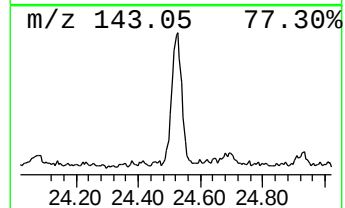
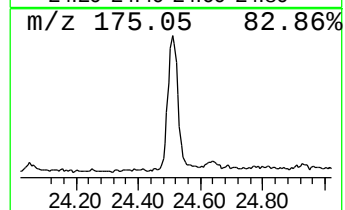
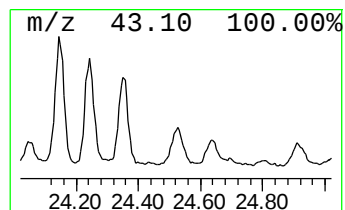
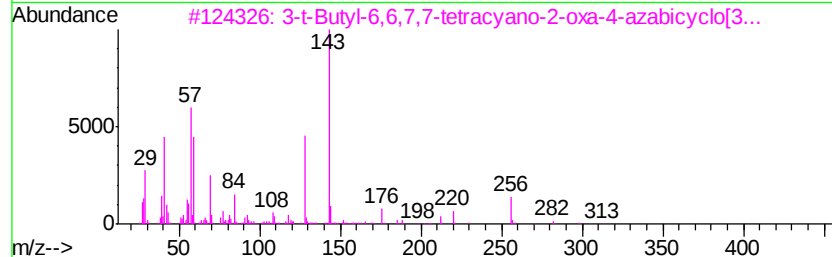
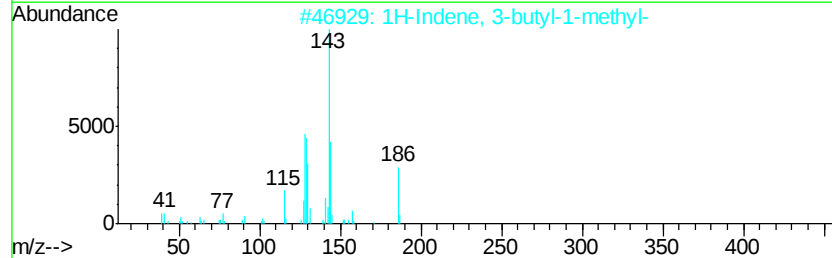
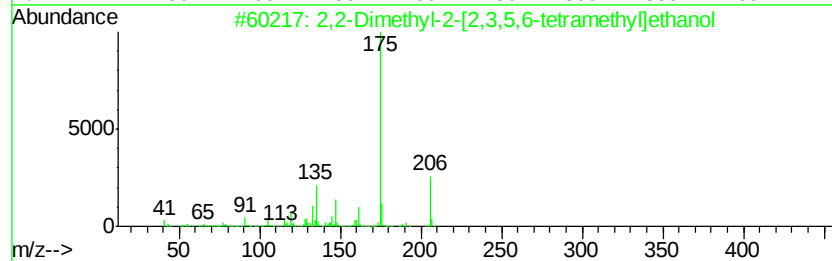
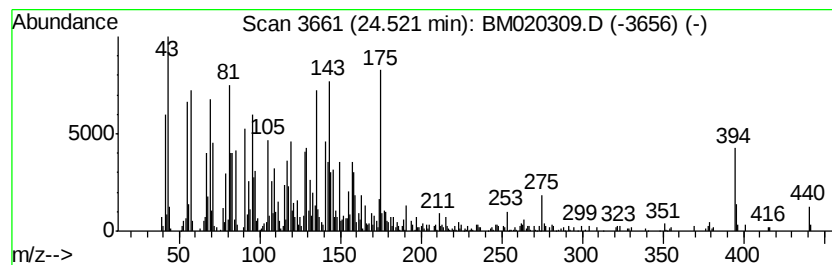
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown24.52 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.52	13.77 ng	1016930	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,2-Dimethyl-2-[2,3,5,6-tetramet...	206	C14H22O	1000128-94-3	16
2		1H-Indene, 3-butyl-1-methyl-	186	C14H18	111400-84-1	12
3		3-t-Butyl-6,6,7,7-tetracyano-2-oxa...	313	C15H15N5O3	1000192-86-4	11
4		Borinic acid, diethyl-, 4-acetyl...	204	C12H17BO2	061142-59-4	11
5		1H-Indole-2-carboxylic acid, 1-m...	175	C10H9NO2	016136-58-6	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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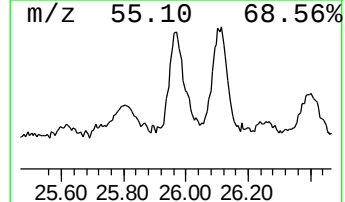
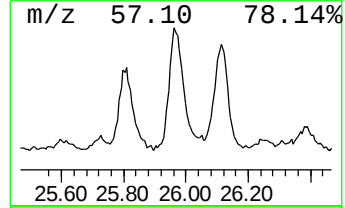
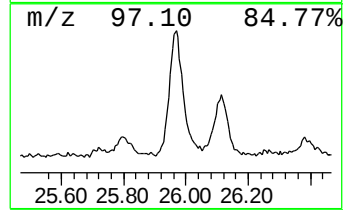
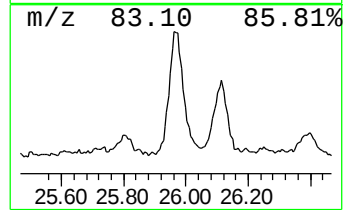
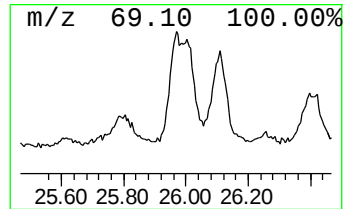
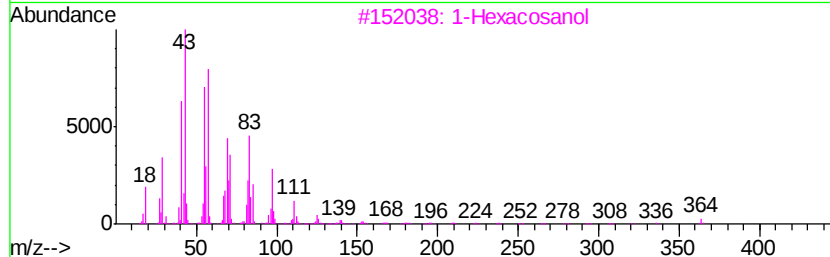
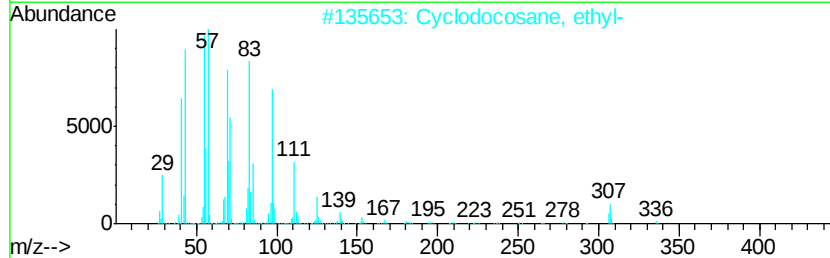
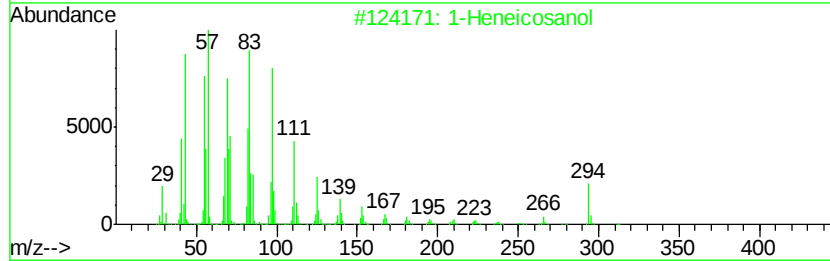
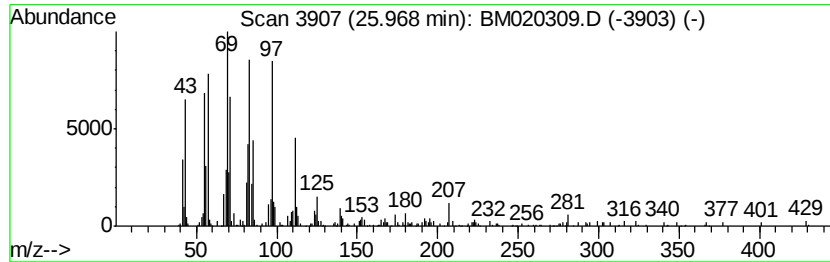
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 1-Heneicosanol Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.97	13.27 ng	979665	Perylene-d12	23.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Heneicosanol	312 C21H44O	015594-90-8	94
2	Cyclodocosane, ethyl-	336 C24H48	1000151-22-6	87
3	1-Hexacosanol	382 C26H54O	000506-52-5	72
4	7-Hexadecene, (Z)-	224 C16H32	035507-09-6	64
5	Cyclohexane, 1,2-dimethyl-3-pent...	224 C16H32	062376-17-4	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
Acq On : 12 May 2019 21:30
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Sample : K2766-10
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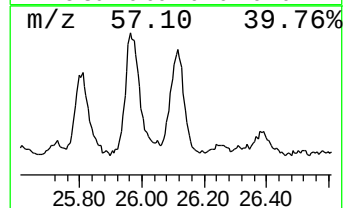
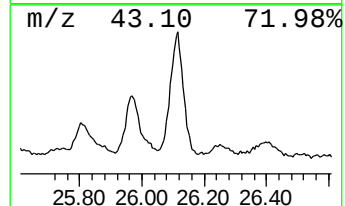
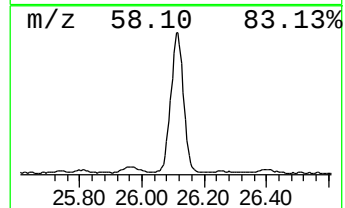
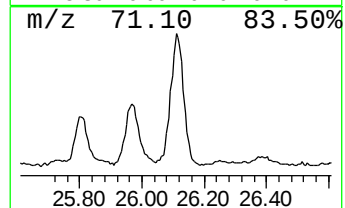
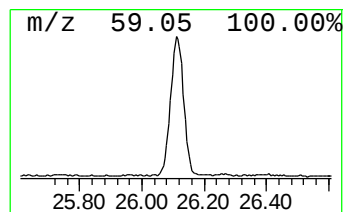
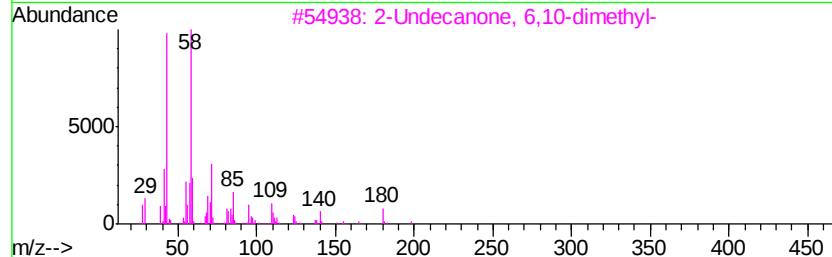
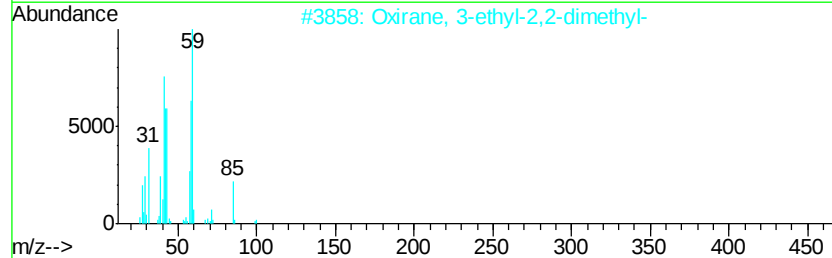
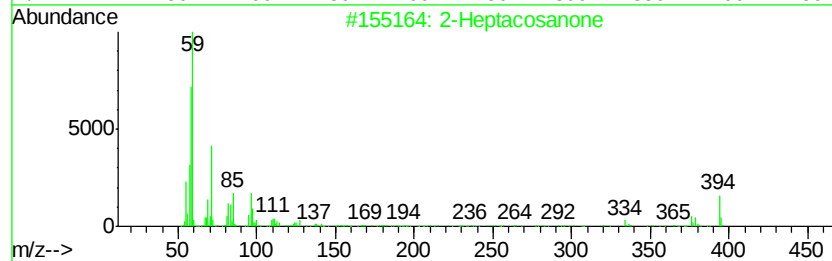
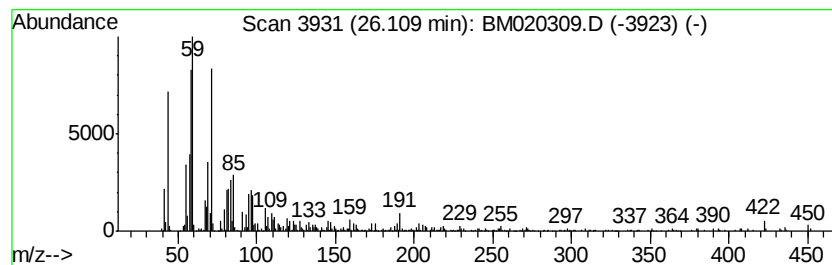
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 2-Heptacosanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.11	26.80 ng	1978960	Perylene-d12	23.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Heptacosanone	394 C27H54O	007796-19-2	59
2	Oxirane, 3-ethyl-2,2-dimethyl-	100 C6H12O	001192-22-9	47
3	2-Undecanone, 6,10-dimethyl-	198 C13H26O	001604-34-8	38
4	2,15-Hexadecanedione	254 C16H30O2	018650-13-0	35
5	Cyclohexanol, 5-methyl-2-(1-meth...	156 C10H20O	000491-02-1	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
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Sample : K2766-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

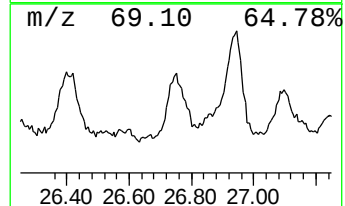
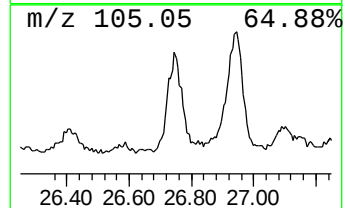
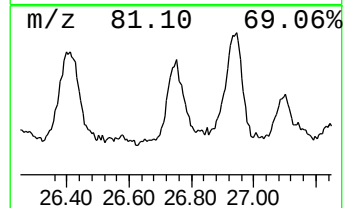
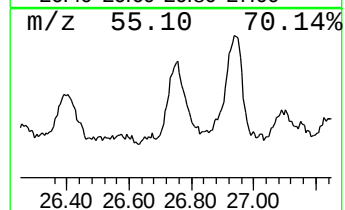
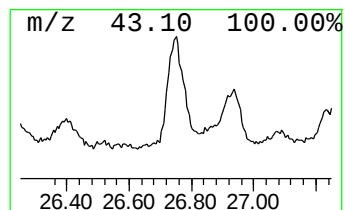
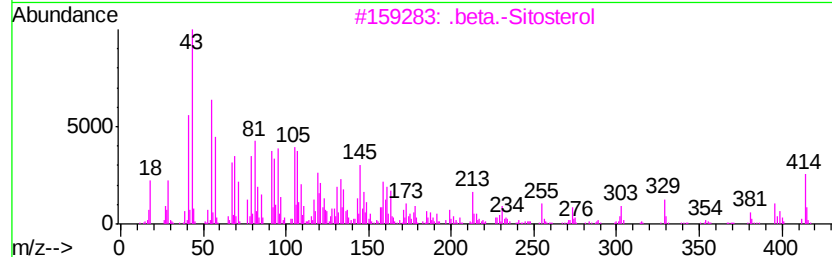
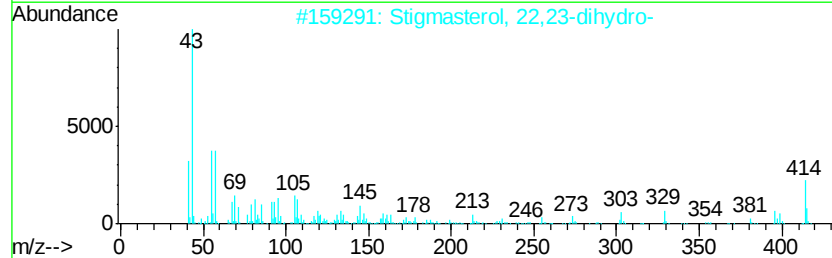
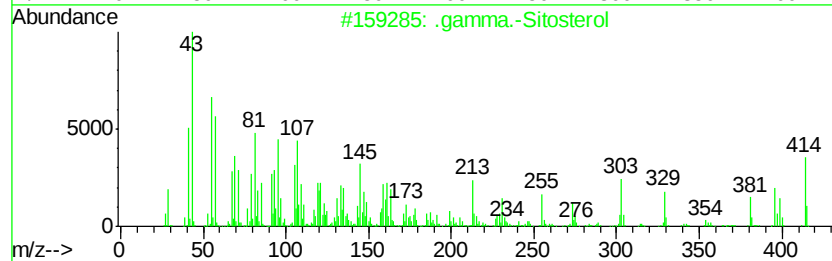
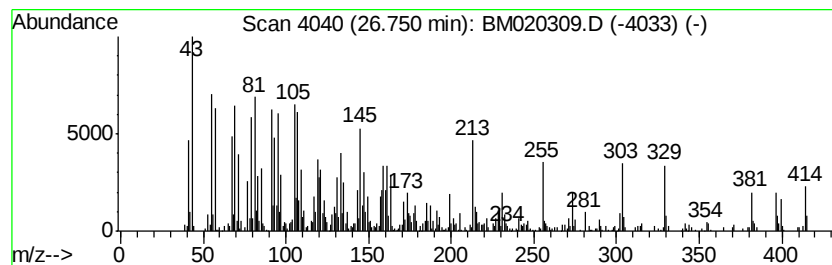
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 .gamma.-Sitosterol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.75	21.01 ng	1551970	Perylene-d12	23.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2	Stigmasterol, 22,23-dihydro-	414	C29H50O	1000214-20-7	99
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	98
4	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	95
5	.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
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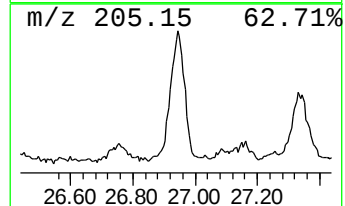
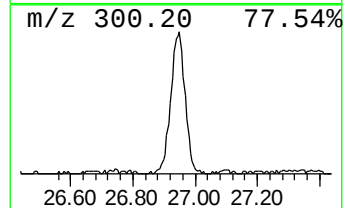
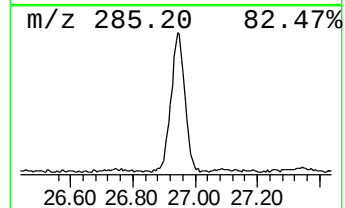
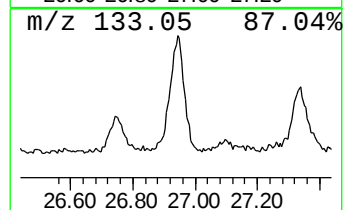
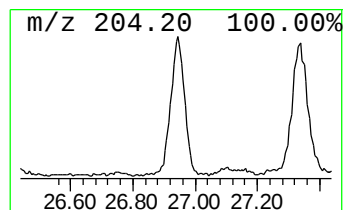
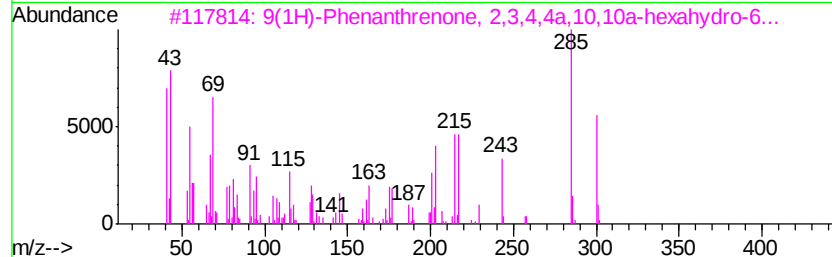
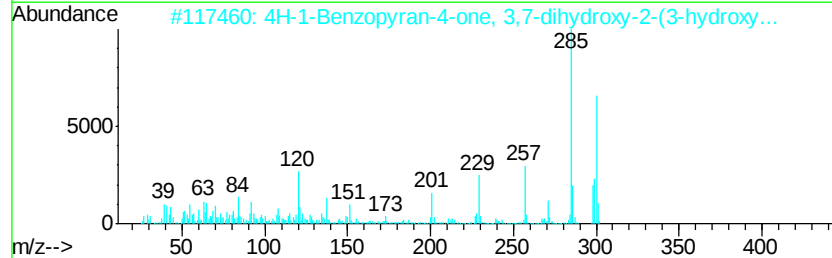
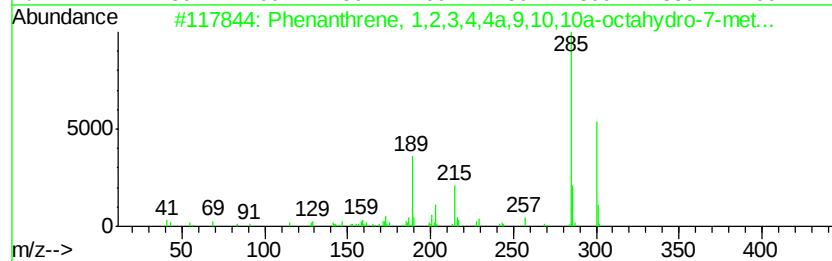
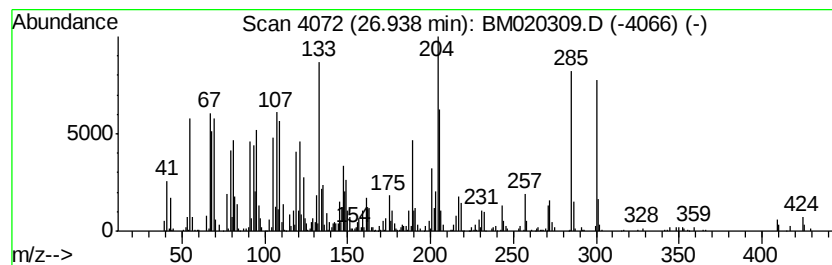
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown26.94 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.94	31.26 ng	2308380	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1,2,3,4,4a,9,10,10...	300	C21H32O	015340-83-7	38
2		4H-1-Benzopyran-4-one, 3,7-dihyd...	300	C16H12O6	057396-72-2	35
3		9(1H)-Phenanthrenone, 2,3,4,4a,1...	300	C20H28O2	000511-05-7	35
4		1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	30
5		1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
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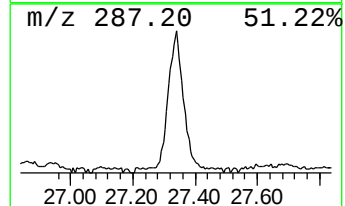
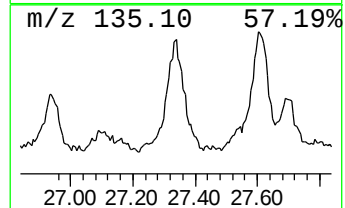
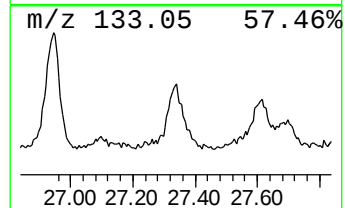
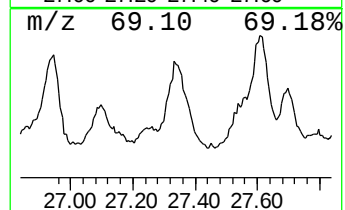
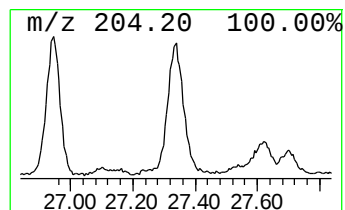
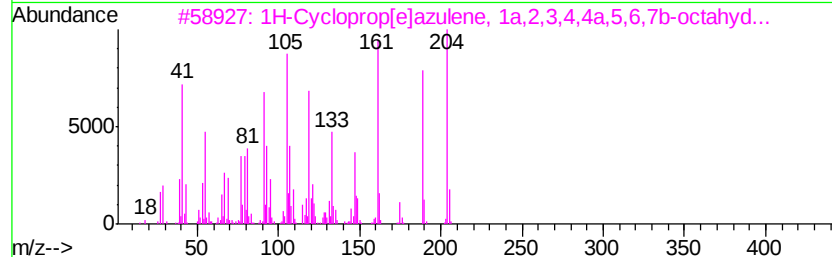
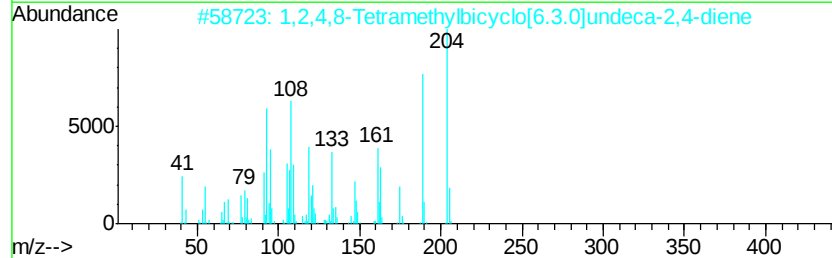
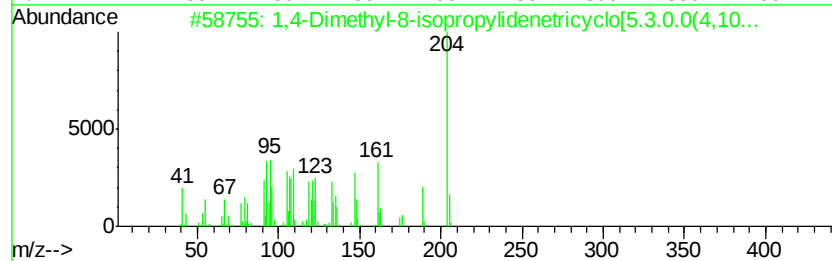
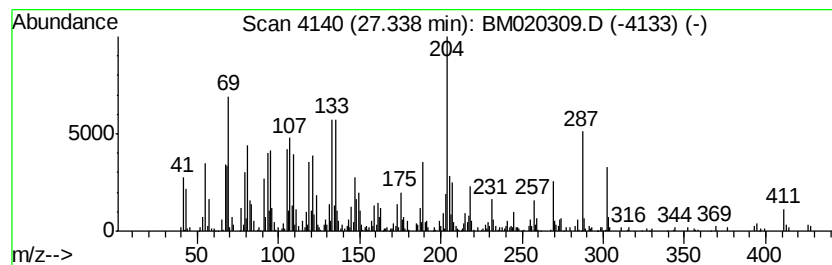
Instrument :
BNA_M
ClientSampleId :
P026-SS008-0002-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 1,4-Dimethyl-8-isopropylide... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.34	30.57 ng	2257630	Perylene-d12	23.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,4-Dimethyl-8-isopropylidenetri...	204 C15H24	1000140-07-7	55
2	1,2,4,8-Tetramethylbicyclo[6.3.0]...	204 C15H24	137235-51-9	43
3	1H-Cycloprop[elazulene, 1a,2,3,4...	204 C15H24	000489-40-7	41
4	1,4-Methanoazulen-9-ol, decahydr...	222 C15H26O	000465-24-7	38
5	2H-Cyclopentacyclooctene, 4,5,6,...	204 C15H24	1000221-85-8	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P026-SS008-0002-01

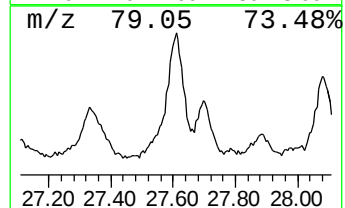
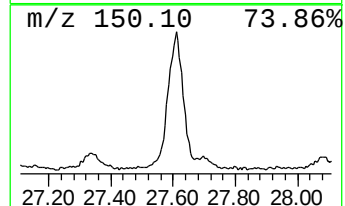
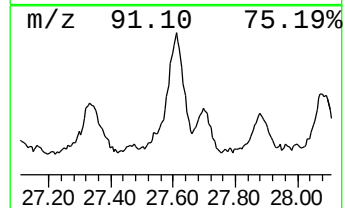
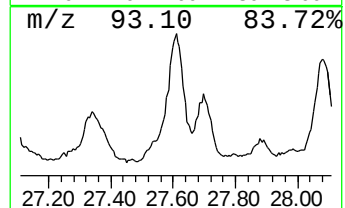
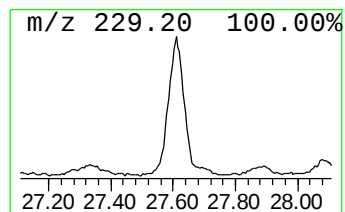
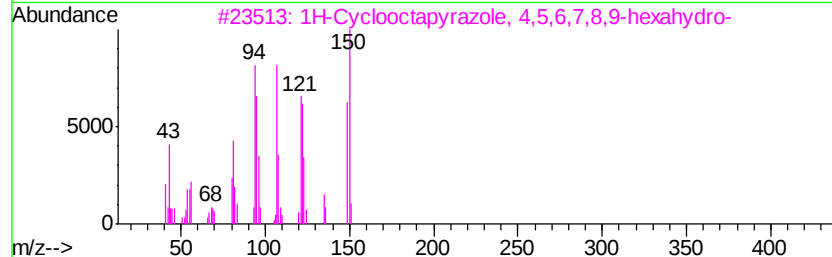
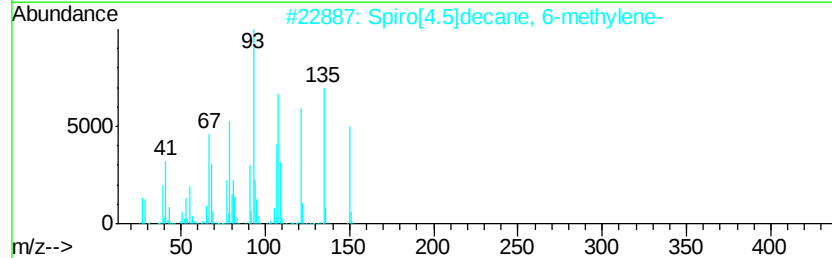
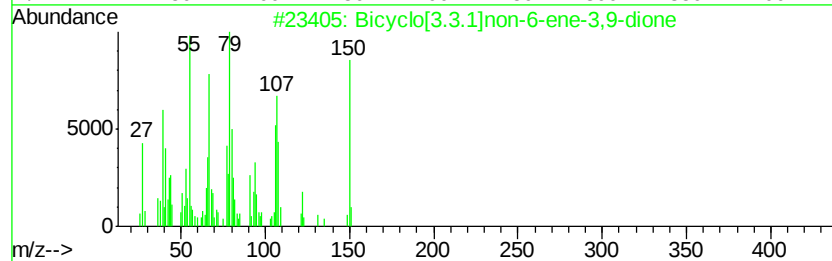
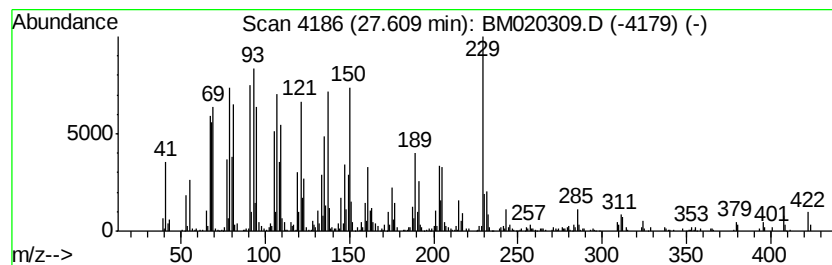
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 unknown27.61 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.61	20.64 ng	1524410	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[3.3.1]non-6-ene-3,9-dione	150	C9H10O2	1000193-44-0	47
2		Spiro[4.5]decane, 6-methylene-	150	C11H18	019144-01-5	42
3		1H-Cyclooctapyrazole, 4,5,6,7,8,...	150	C9H14N2	015984-10-8	38
4		Butyramide, 3-(2-furyl)-N-phenyl-	229	C14H15NO2	1000156-94-9	38
5		Cyclohexene, 2,4-dimethyl-1-(1-m...	150	C11H18	056763-60-1	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051219\
Data File : BM020309.D
Acq On : 12 May 2019 21:30
Operator : JU/SJ
Sample : K2766-10
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P026-SS008-0002-01

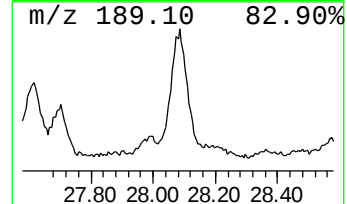
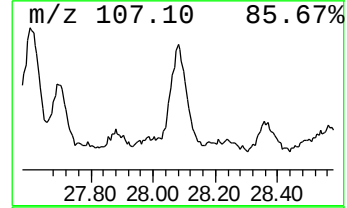
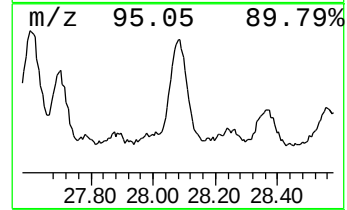
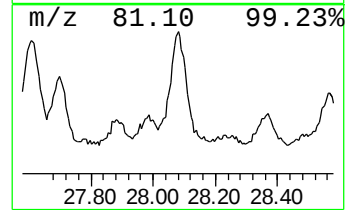
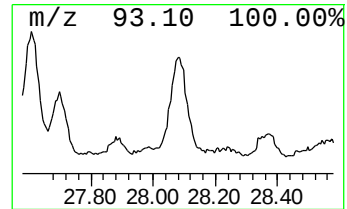
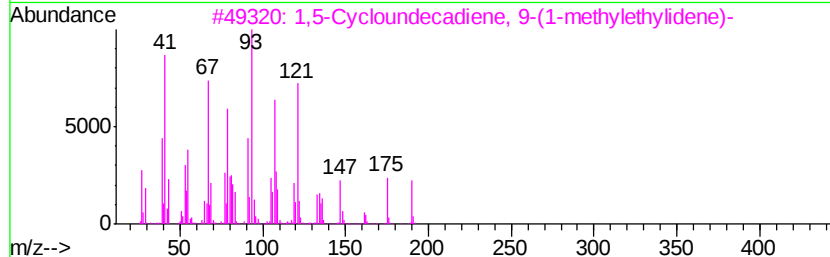
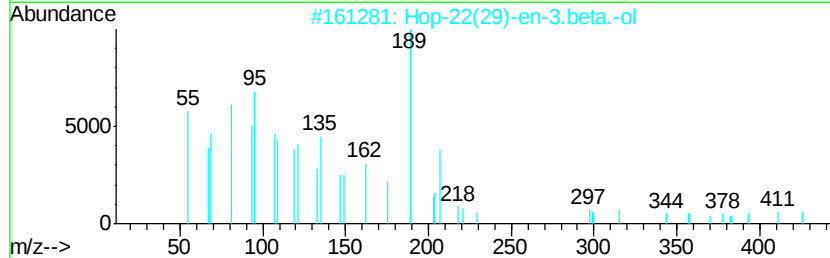
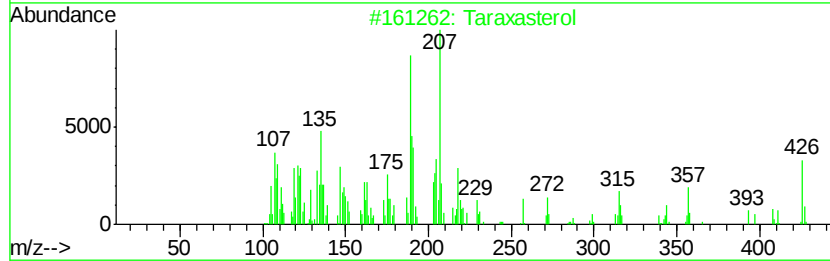
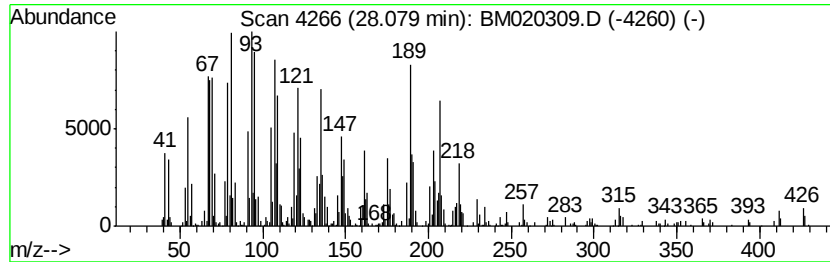
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Taraxasterol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.08	29.91 ng	2209010	Perylene-d12	23.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Taraxasterol	426	C30H50O	001059-14-9	60
2		Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	50
3		1,5-Cycloundecadiene, 9-(1-methv...	190	C14H22	062338-55-0	40
4		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	25
5		1,5,9-Cyclododecatriene, 1,5,9-t...	204	C15H24	021064-19-7	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM051219\
 Data File : BM020309.D
 Acq On : 12 May 2019 21:30
 Operator : JU/SJ
 Sample : K2766-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P026-SS008-0002-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM043019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Toluene	3.98	13.3	ng	307569	1	7.76	463317	20.0
unknown7.29	7.29	80.1	ng	1855900	1	7.76	463317	20.0
n-Hexadecanoic acid	18.04	21.3	ng	1343070	4	17.16	1263340	20.0
1,1'-Biphenyl, 2,...	20.06	12.3	ng	1895520	5	21.35	3085500	20.0
Ethanol, 2-(tetra...	21.04	31.1	ng	4796120	5	21.35	3085500	20.0
3-Eicosene, (E)-	21.94	63.4	ng	9787190	5	21.35	3085500	20.0
Heneicosane	22.91	28.5	ng	2104120	6	23.64	1477040	20.0
1-Heneicosyl formate	22.96	42.9	ng	3170780	6	23.64	1477040	20.0
Hexadecane, 1-iodo-	24.14	20.2	ng	1491360	6	23.64	1477040	20.0
1-Hexacosanol	24.24	25.8	ng	1904040	6	23.64	1477040	20.0
2-Nonacosanone	24.35	13.1	ng	968976	6	23.64	1477040	20.0
unknown24.52	24.52	13.8	ng	1016930	6	23.64	1477040	20.0
1-Heneicosanol	25.97	13.3	ng	979665	6	23.64	1477040	20.0
2-Heptacosanone	26.11	26.8	ng	1978960	6	23.64	1477040	20.0
.gamma.-Sitosterol	26.75	21.0	ng	1551970	6	23.64	1477040	20.0
unknown26.94	26.94	31.3	ng	2308380	6	23.64	1477040	20.0
1,4-Dimethyl-8-is...	27.34	30.6	ng	2257630	6	23.64	1477040	20.0
unknown27.61	27.61	20.6	ng	1524410	6	23.64	1477040	20.0
Taraxasterol	28.08	29.9	ng	2209010	6	23.64	1477040	20.0