

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
 Data File : BF120957.D
 Acq On : 21 Jul 2020 16:46
 Operator : JU/CG
 Sample : L3354-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BACK-YARD

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_F\METHODS\8270-BF071620.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	5.422	563	567	571	rBV	1880752	1792367	45.26%	5.866%
2	6.439	735	740	743	rBV	2109117	1982740	50.07%	6.489%
3	6.645	772	775	789	rVB	31893	45017	1.14%	0.147%
4	6.816	800	804	807	rBV	1072665	976519	24.66%	3.196%
5	7.375	894	899	902	rBV	1405408	1234083	31.16%	4.039%
6	7.839	975	978	982	rVB	241037	218677	5.52%	0.716%
7	8.098	1017	1022	1025	rBV	1460653	1304299	32.94%	4.269%
8	8.716	1122	1127	1131	rBV	2887355	2810302	70.97%	9.197%
9	9.175	1200	1205	1208	rBV	2914569	2777204	70.13%	9.089%
10	9.269	1218	1221	1223	rBV	133201	104105	2.63%	0.341%
11	9.563	1268	1271	1284	rBV	449837	412570	10.42%	1.350%
12	9.851	1315	1320	1330	rBV	1715055	1565906	39.54%	5.125%
13	10.639	1449	1454	1464	rBV	2132097	2285820	57.72%	7.481%
14	11.210	1549	1551	1561	rVB4	33519	55186	1.39%	0.181%
15	11.333	1568	1572	1575	rBV	1944214	1648445	41.63%	5.395%
16	11.868	1659	1663	1666	rBV2	51254	52178	1.32%	0.171%
17	12.210	1716	1721	1728	rBV2	49183	56591	1.43%	0.185%
18	12.310	1732	1738	1741	rBV	259926	231473	5.85%	0.758%
19	12.368	1745	1748	1751	rVB	46970	39799	1.01%	0.130%
20	12.539	1773	1777	1779	rBV	46294	40808	1.03%	0.134%
21	12.563	1779	1781	1784	rVB	62337	55589	1.40%	0.182%
22	12.768	1810	1816	1819	rBV	52618	79311	2.00%	0.260%
23	12.921	1837	1842	1845	rBV	3975547	3959961	100.00%	12.960%
24	13.151	1876	1881	1884	rBV	72347	84108	2.12%	0.275%
25	13.445	1927	1931	1935	rBV	102469	92728	2.34%	0.303%
26	13.533	1943	1946	1950	rBV	167189	137036	3.46%	0.448%
27	13.768	1982	1986	1991	rVB	99102	98326	2.48%	0.322%
28	13.827	1992	1996	2004	rBV	595315	813897	20.55%	2.664%
29	13.968	2015	2020	2023	rBV	1726396	1514024	38.23%	4.955%
30	14.445	2096	2101	2102	rBV2	101438	108754	2.75%	0.356%
31	14.468	2102	2105	2114	rVB2	119470	192177	4.85%	0.629%
32	14.615	2126	2130	2136	rVB	117058	120893	3.05%	0.396%
33	14.745	2149	2152	2155	rVB	53049	50001	1.26%	0.164%
34	14.892	2173	2177	2182	rVB	161130	161969	4.09%	0.530%

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Instrument :
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.080	2204	2209	2212	rBV	276204	299747	7.57%	0.981%
36	15.121	2212	2216	2229	rVB2	210087	380001	9.60%	1.244%
37	15.415	2261	2266	2270	rBV	1007589	1182154	29.85%	3.869%
38	15.451	2270	2272	2280	rVB	54976	70617	1.78%	0.231%
39	15.651	2302	2306	2314	rVB2	45157	63273	1.60%	0.207%
40	15.880	2340	2345	2351	rBV	303151	424716	10.73%	1.390%
41	15.956	2352	2358	2368	rBV2	80867	221802	5.60%	0.726%
42	16.374	2425	2429	2433	rBV2	26202	40784	1.03%	0.133%
43	16.656	2471	2477	2483	rBV4	41264	77271	1.95%	0.253%
44	16.951	2522	2527	2537	rVB2	57182	118176	2.98%	0.387%
45	17.074	2539	2548	2554	rBV8	48483	152775	3.86%	0.500%
46	17.186	2563	2567	2576	rVB5	43719	90417	2.28%	0.296%
47	17.445	2604	2611	2622	rBV6	72122	198047	5.00%	0.648%
48	17.909	2683	2690	2697	rBV10	26312	65657	1.66%	0.215%
49	18.427	2771	2778	2790	rVB10	19636	67759	1.71%	0.222%

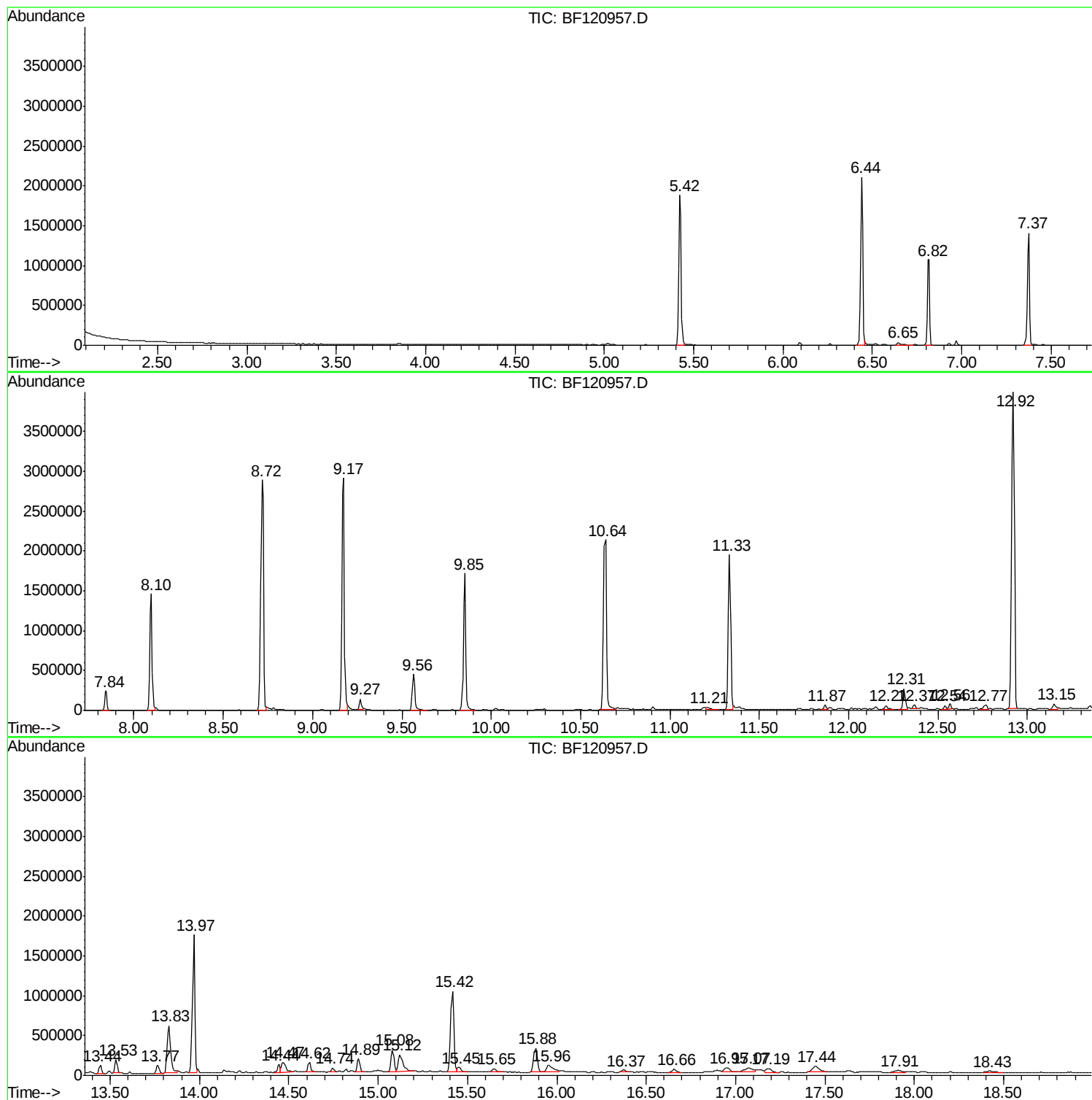
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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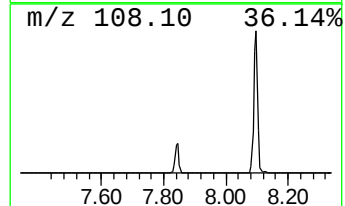
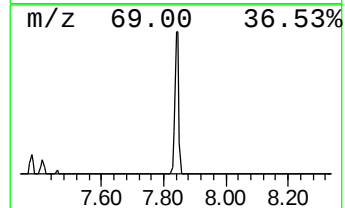
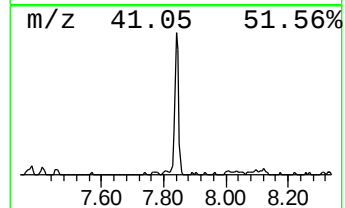
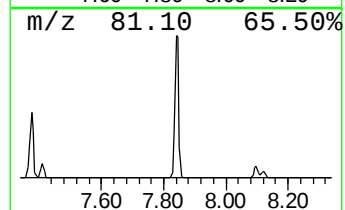
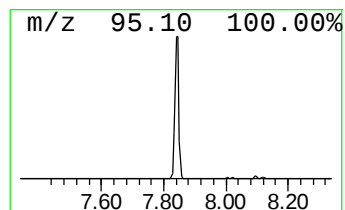
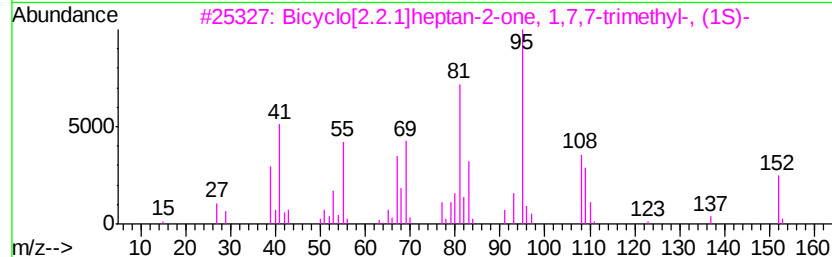
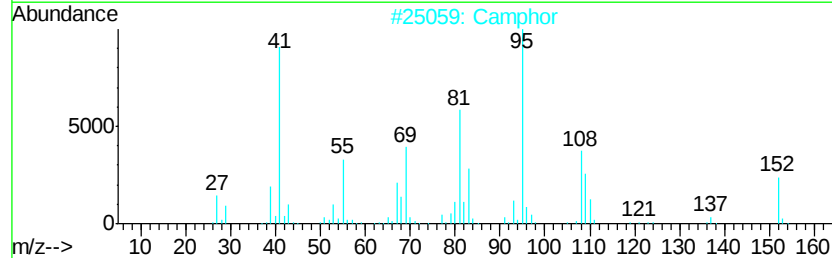
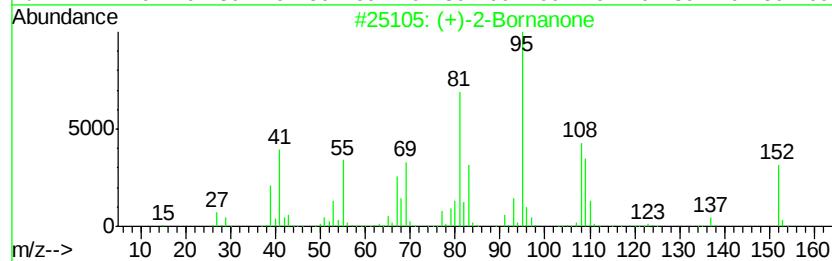
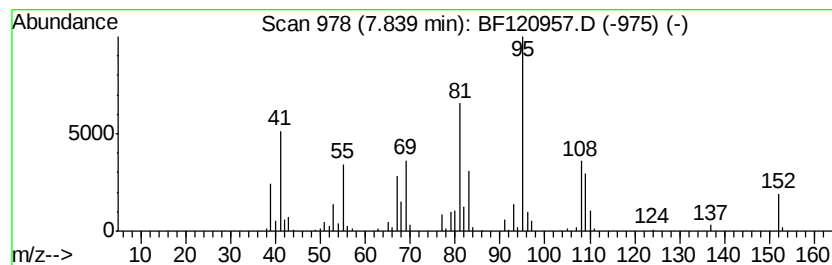
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 (+)-2-Bornanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.84	3.35 ng	218677	Napthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(+)-2-Bornanone	152	C10H16O	000464-49-3	97
2		Camphor	152	C10H16O	000076-22-2	97
3		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	96
4		5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	006752-80-3	58
5		1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	46



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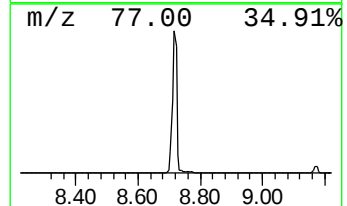
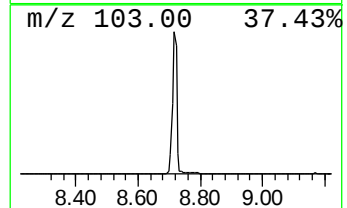
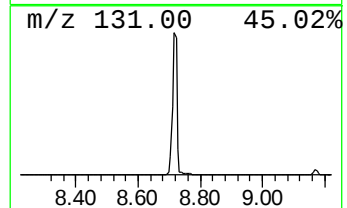
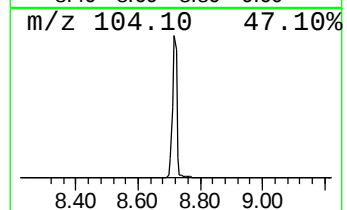
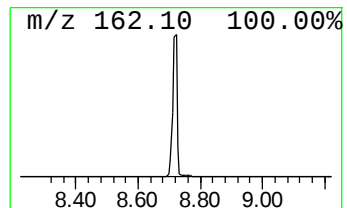
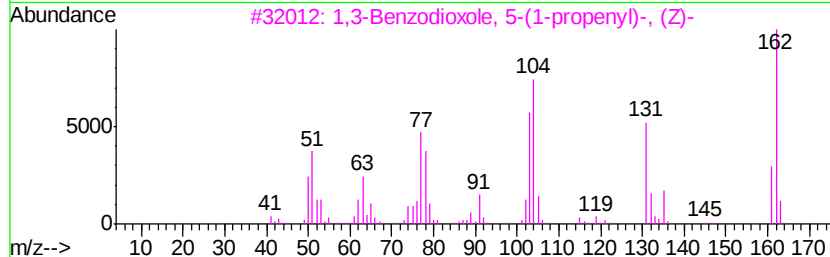
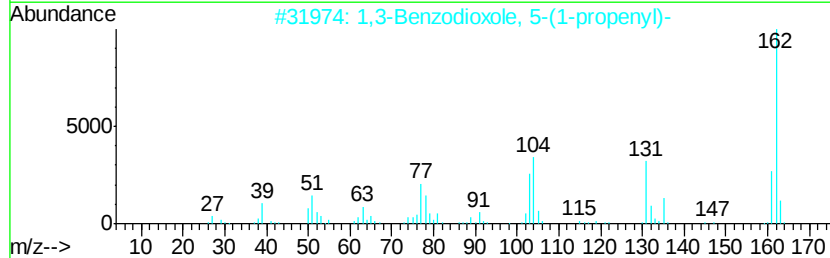
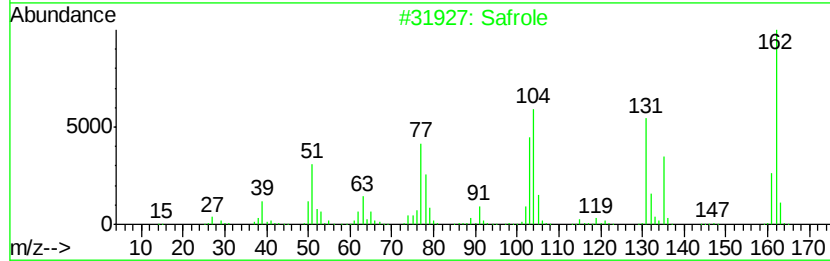
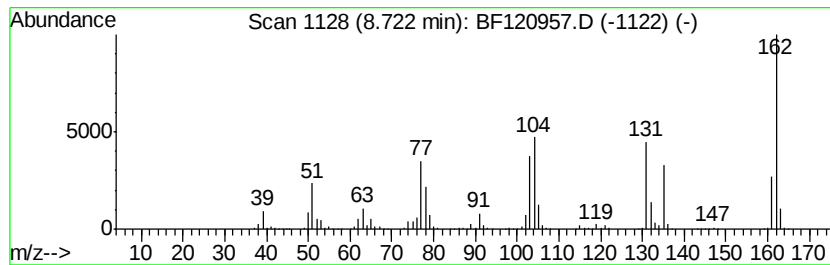
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Safrole Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	43.09 ng	2810300	Naphthalene-d8	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Safrole	162	C10H10O2	000094-59-7	98
2		1,3-Benzodioxole, 5-(1-propenyl)-	162	C10H10O2	000120-58-1	97
3		1,3-Benzodioxole, 5-(1-propenyl)...	162	C10H10O2	017627-76-8	96
4		Benzene, 1,2-(methylenedioxy)-4-...	162	C10H10O2	004043-71-4	95
5		1,4-Dioxo-1,2,3,4-tetrahydrophth...	162	C8H6N2O2	001445-69-8	41



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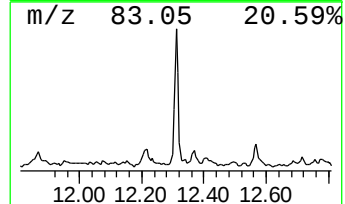
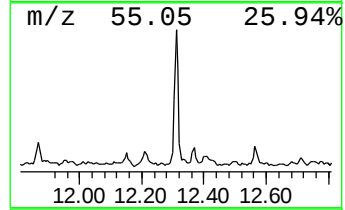
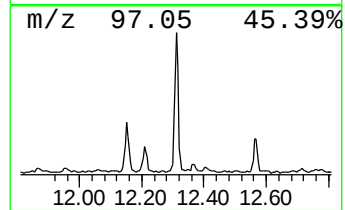
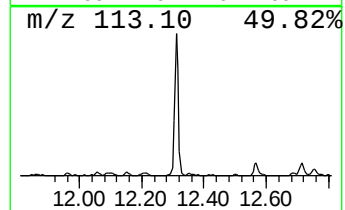
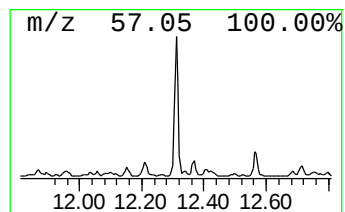
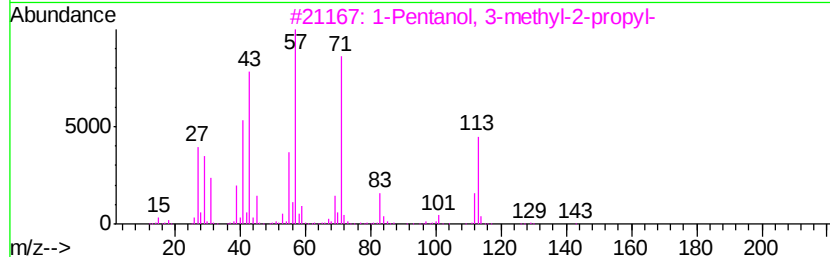
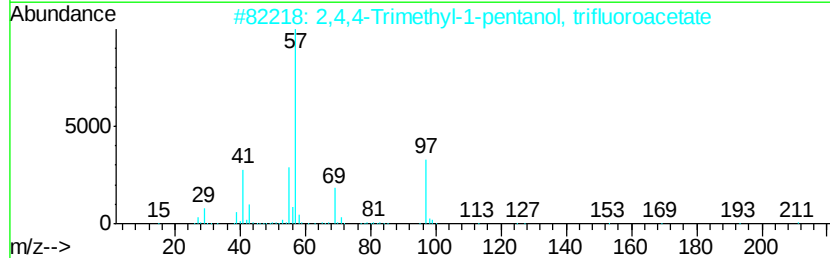
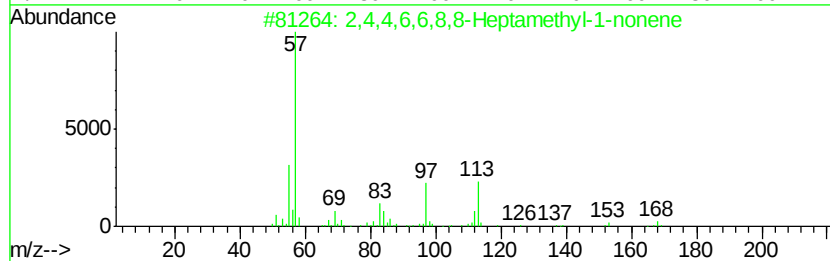
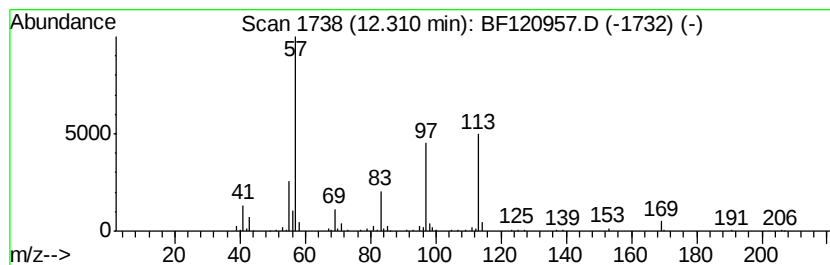
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2,4,4,6,6,8,8-Heptamethyl-1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.31	2.81 ng	231473	Phenanthrene-d10	11.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	56		
2	2,4,4-Trimethyl-1-pentanol, trif...	226	C10H17F3O2	1000365-19-5	35		
3	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	33		
4	Sulfurous acid, di(2-ethylhexyl)...	306	C16H34O3S	1000309-19-1	32		
5	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	32		



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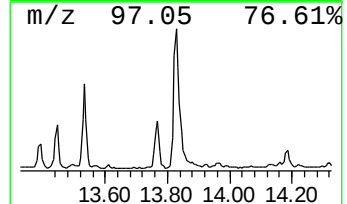
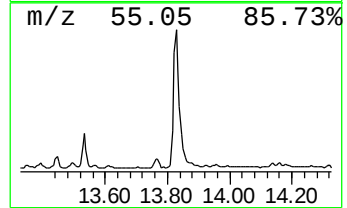
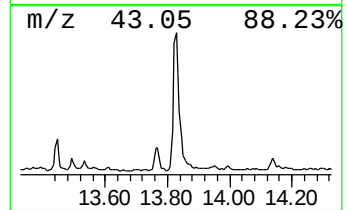
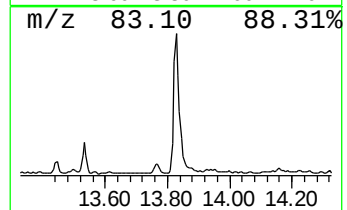
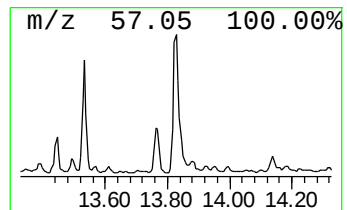
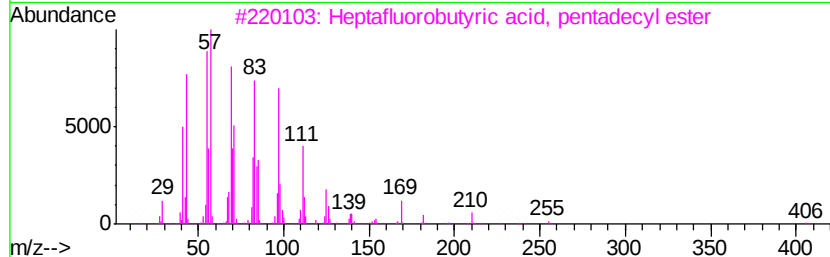
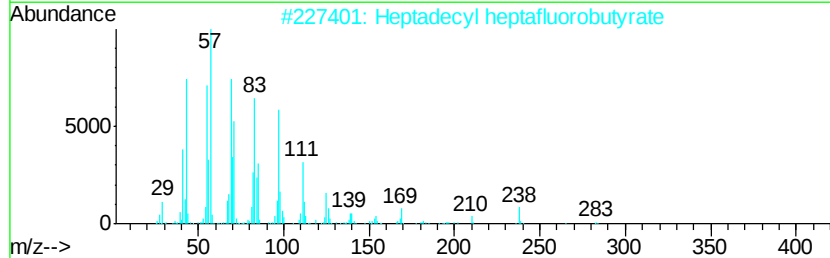
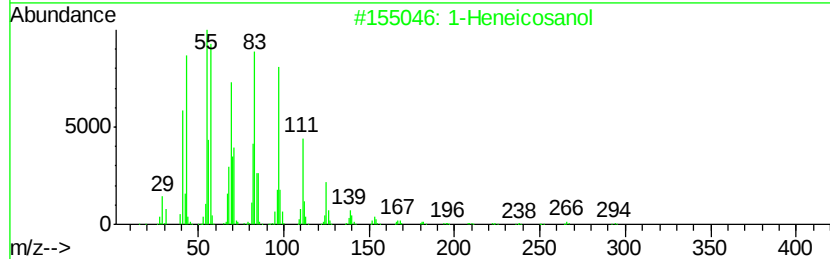
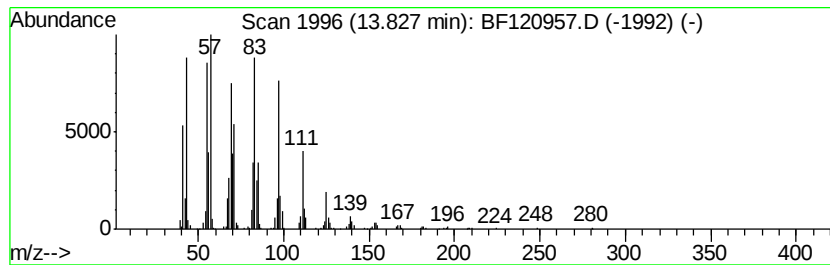
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	10.75 ng	813897	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	95
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4		Behenic alcohol	326	C22H46O	000661-19-8	94
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	93



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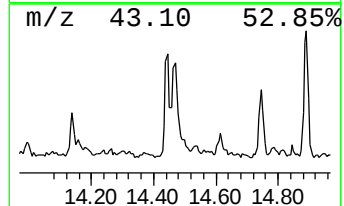
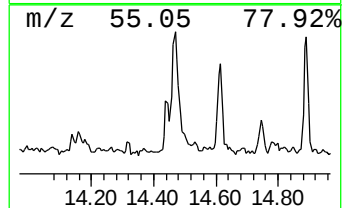
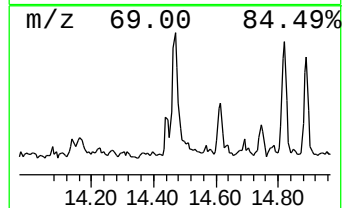
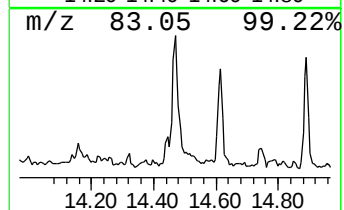
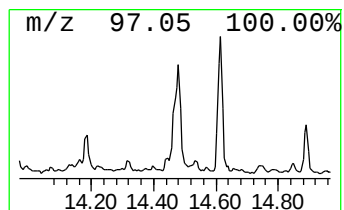
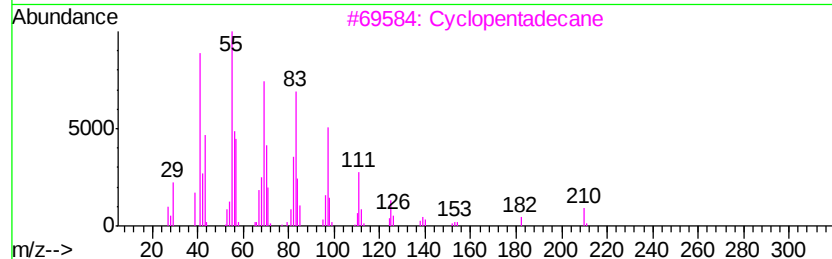
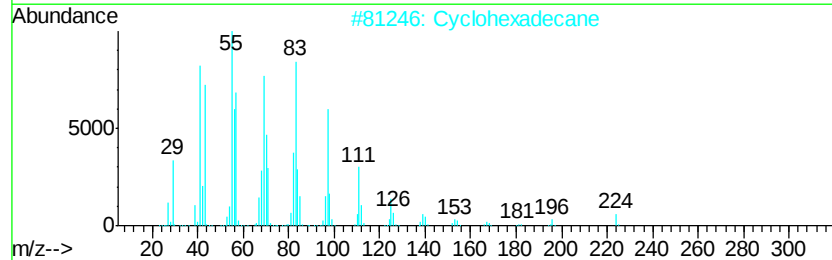
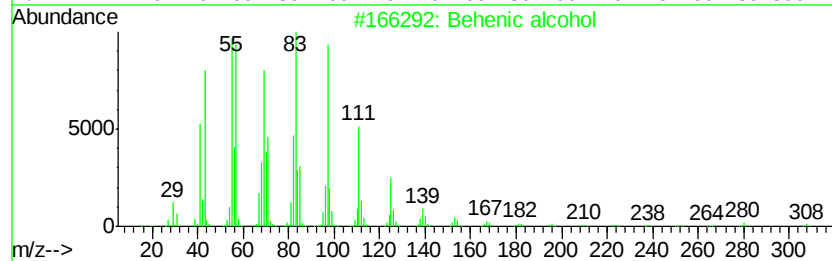
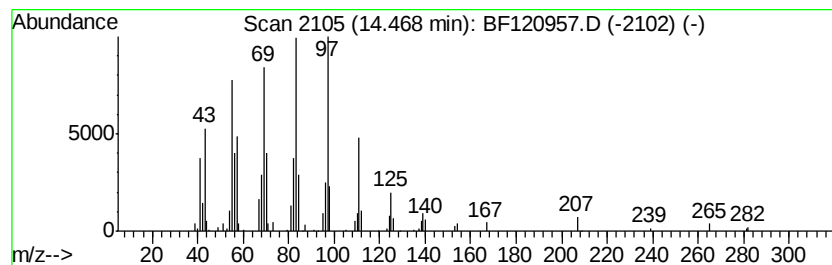
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Cyclopentadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.47	2.54 ng	192177	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	94
2		Cyclohexadecane	224	C16H32	000295-65-8	90
3		Cyclopentadecane	210	C15H30	000295-48-7	86
4		1-Heneicosanol	312	C21H44O	015594-90-8	83
5		n-Pentadecanol	228	C15H32O	000629-76-5	72



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Acq On : 21 Jul 2020 16:46
Operator : JU/CG
Sample : L3354-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

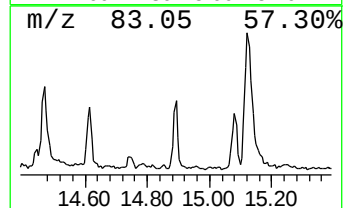
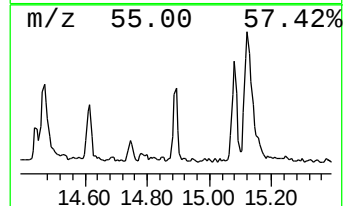
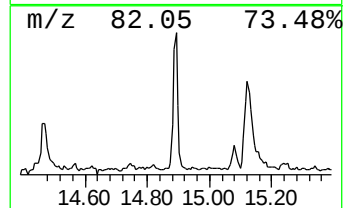
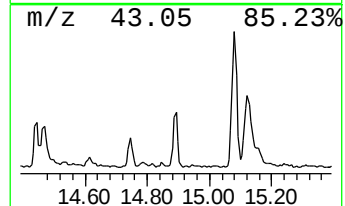
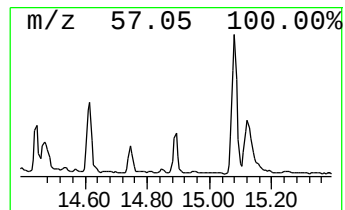
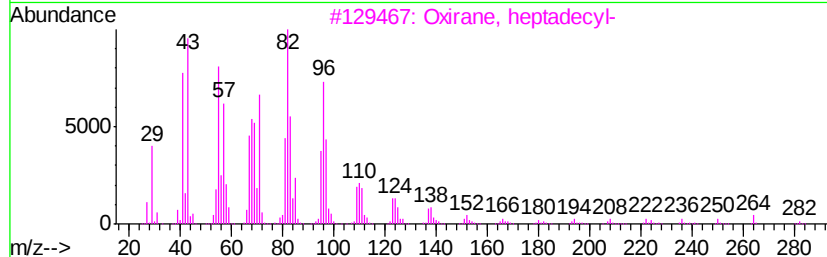
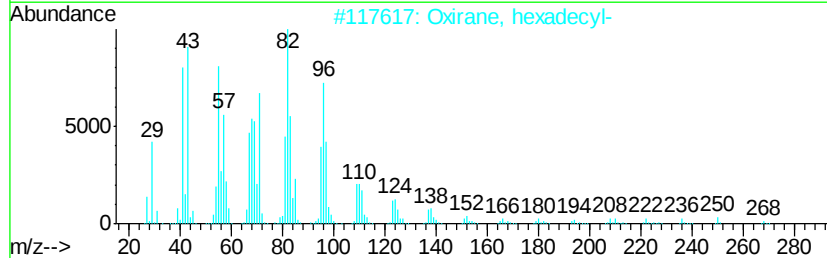
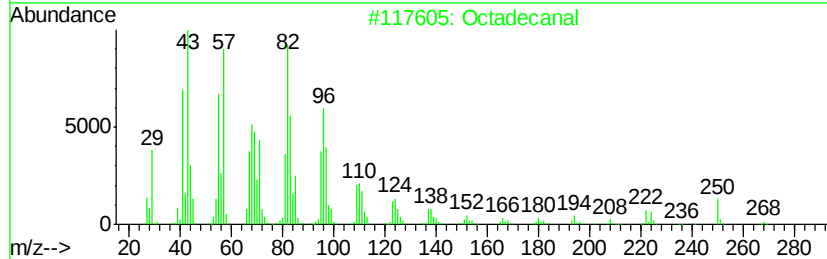
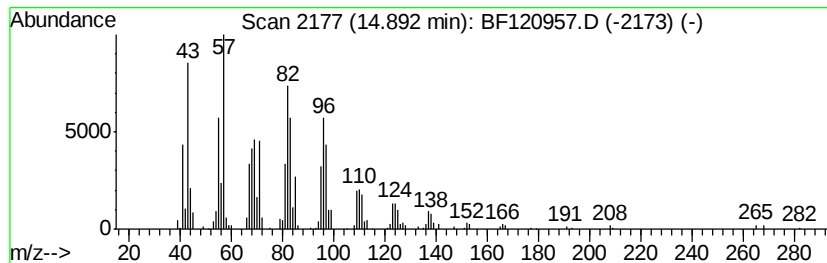
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.89	2.74 ng	161969	Perylene-d12	15.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanal	268	C18H36O	000638-66-4	99
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	90
4	1,19-Eicosadiene	278	C20H38	014811-95-1	87
5	Oleyl alcohol, methyl ether	282	C19H38O	1000352-68-0	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
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ClientSampleId :
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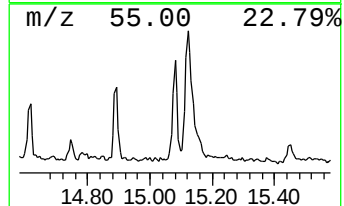
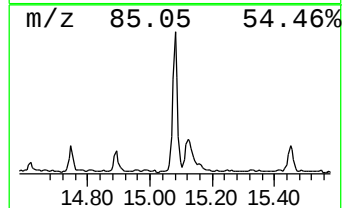
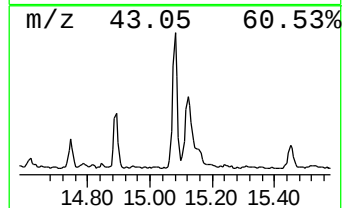
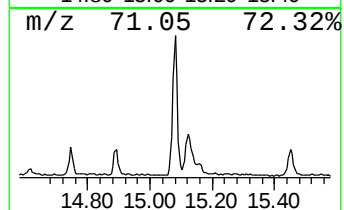
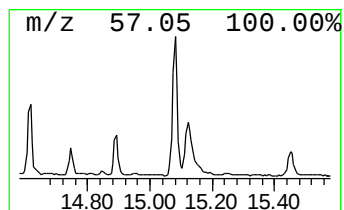
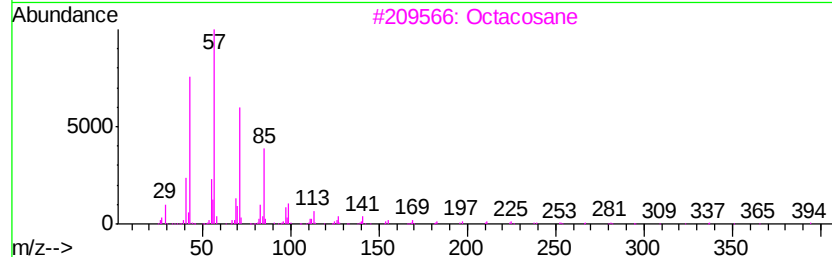
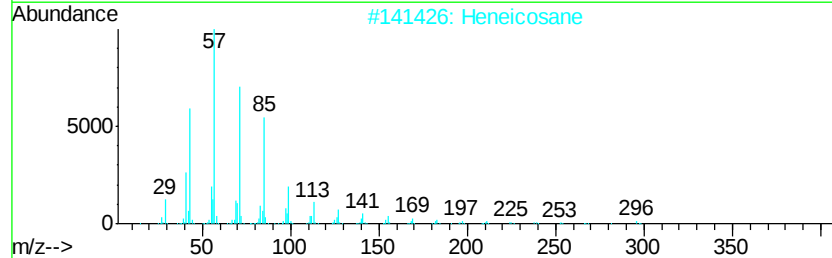
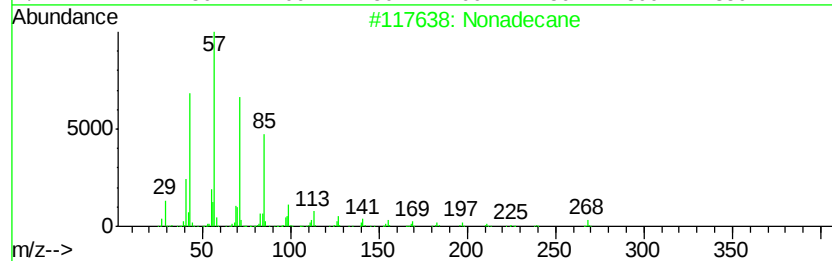
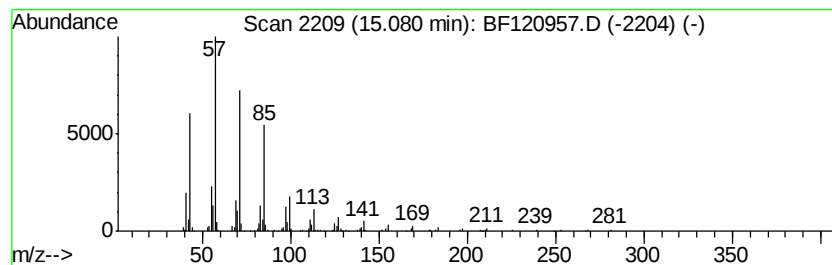
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	5.07 ng	299747	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	94
2	Heneicosane		296	C21H44	000629-94-7	91
3	Octacosane		394	C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
5	Hexacosane		366	C26H54	000630-01-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
Data File : BF120957.D
Acq On : 21 Jul 2020 16:46
Operator : JU/CG
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Misc :
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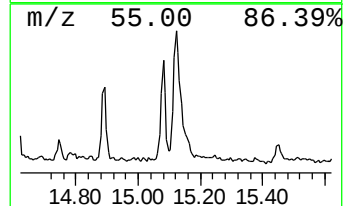
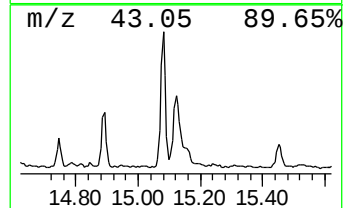
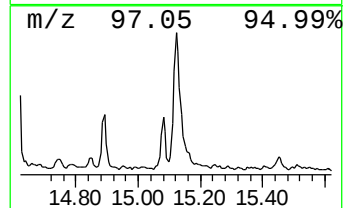
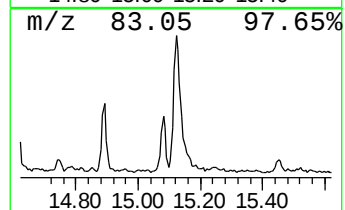
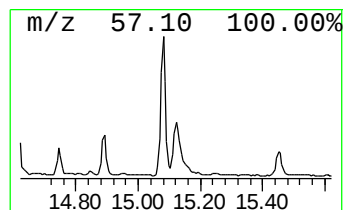
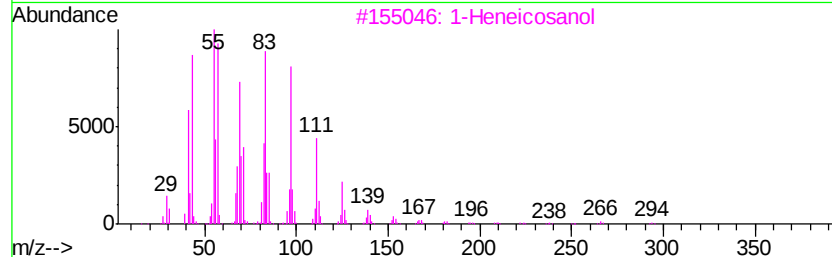
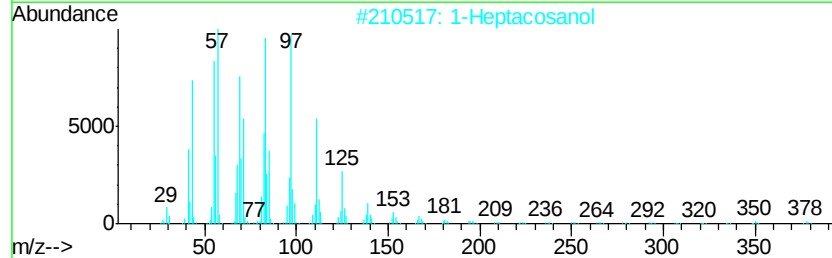
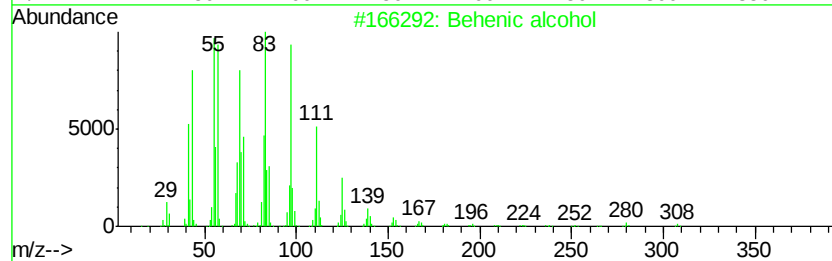
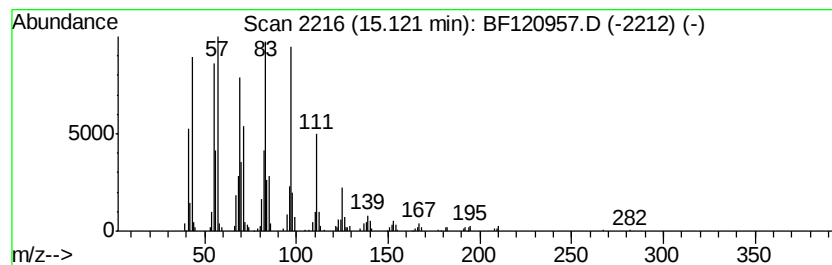
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Behenic alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	6.43 ng	380001	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Behenic alcohol	326	C22H46O	000661-19-8	95
2		1-Heptacosanol	396	C27H56O	002004-39-9	95
3		1-Heneicosanol	312	C21H44O	015594-90-8	95
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
5		Octacosanol	410	C28H58O	000557-61-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
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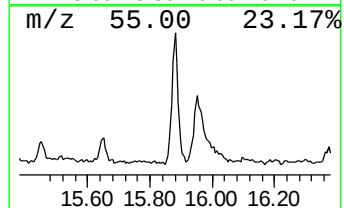
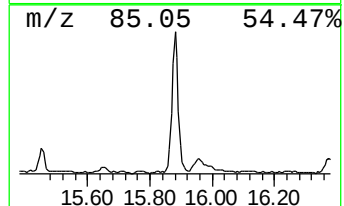
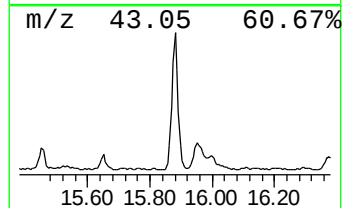
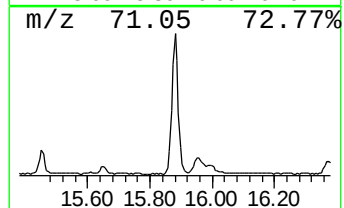
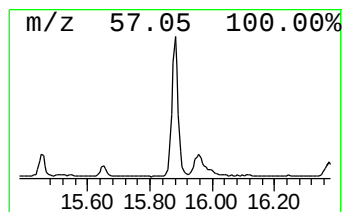
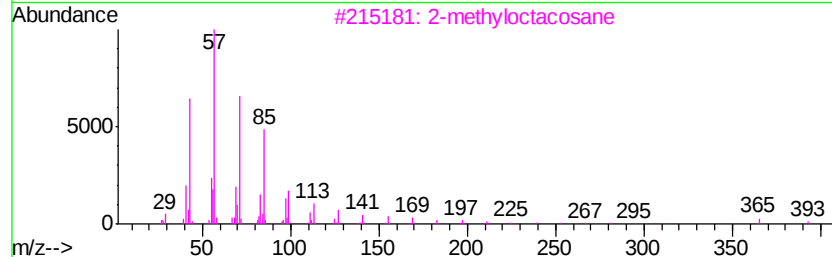
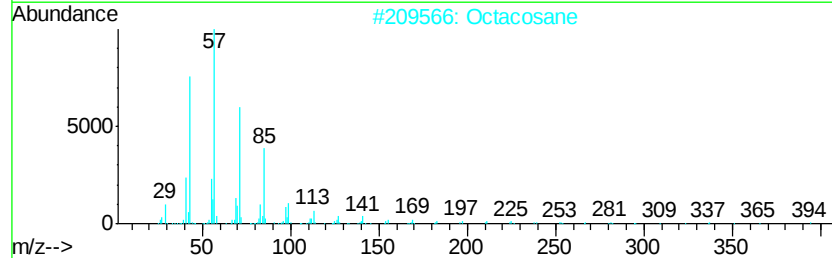
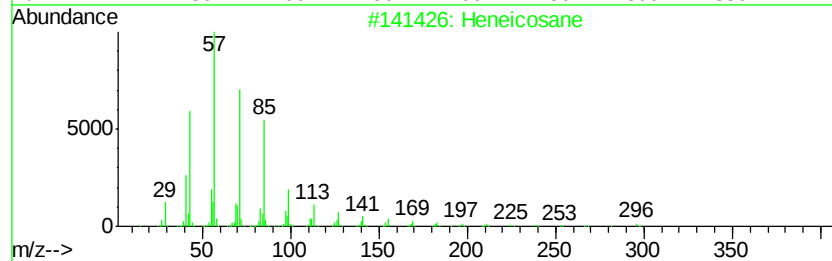
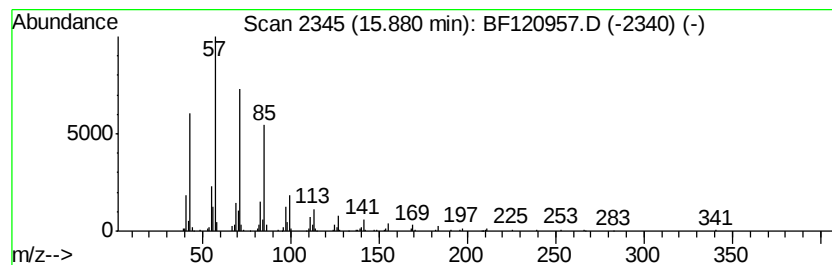
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Heneicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	7.19 ng	424716	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Octacosane		394	C28H58	000630-02-4	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Hexacosane		366	C26H54	000630-01-3	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
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Misc :
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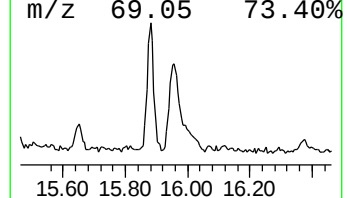
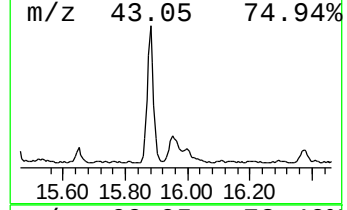
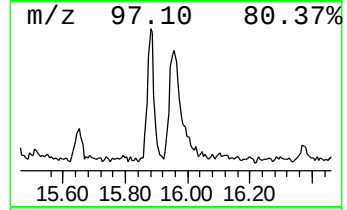
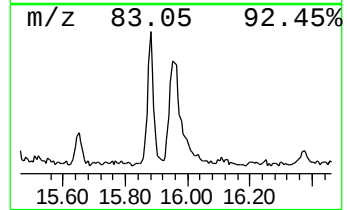
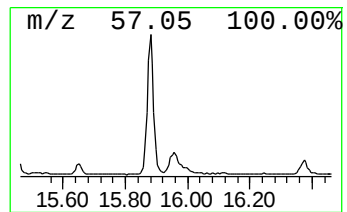
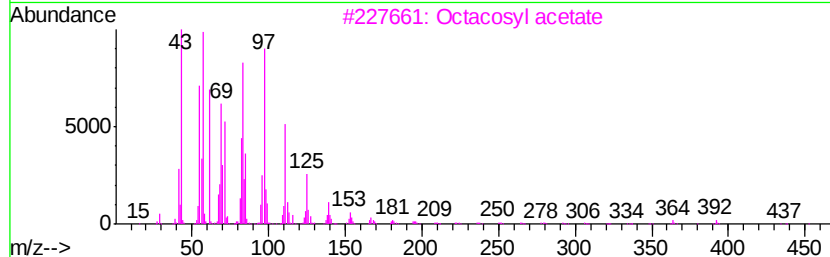
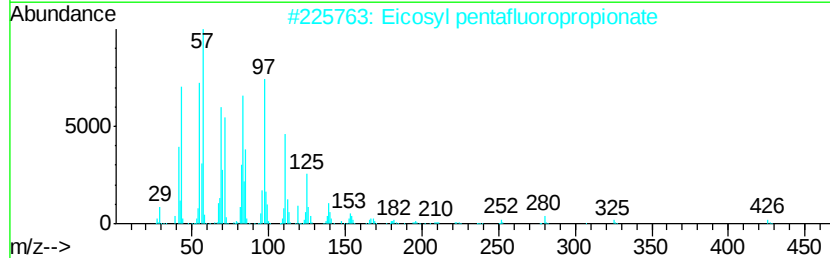
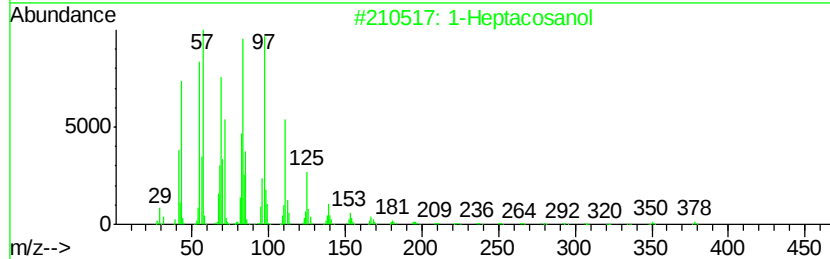
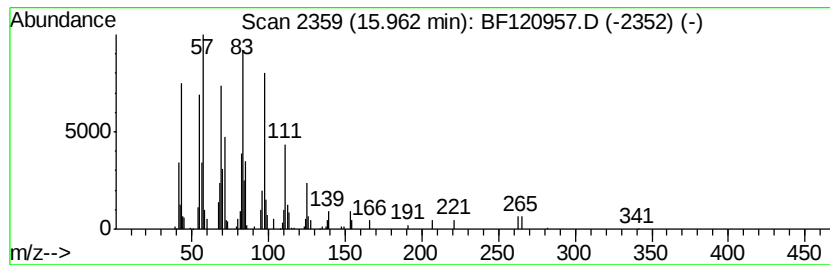
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Heptacosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	3.75 ng	221802	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Heptacosanol	396 C27H56O	002004-39-9	95
2	Eicosyl pentafluoropropionate	444 C23H41F5O2	1000351-80-8	94
3	Octacosyl acetate	452 C30H60O2	018206-97-8	94
4	Heptafluorobutyric acid, n-octad...	466 C22H37F7O2	000400-57-7	94
5	n-Tetracosanol-1	354 C24H50O	000506-51-4	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
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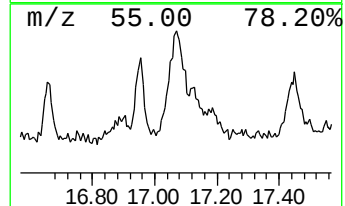
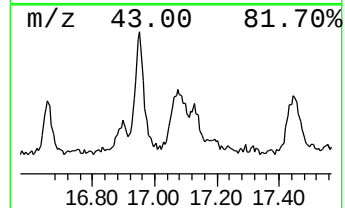
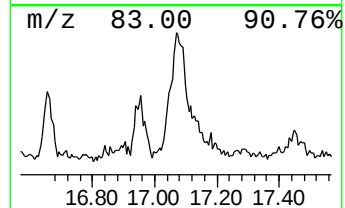
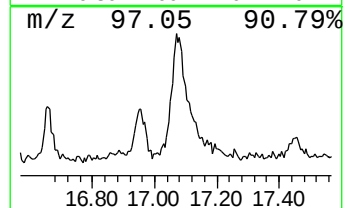
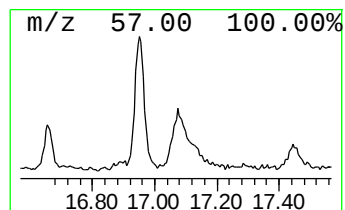
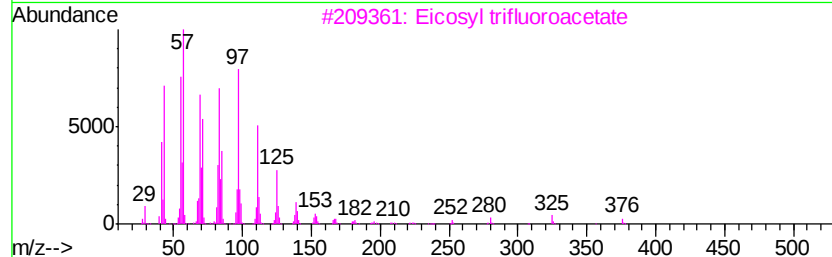
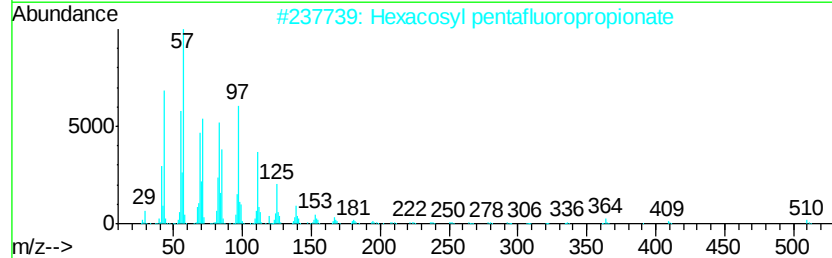
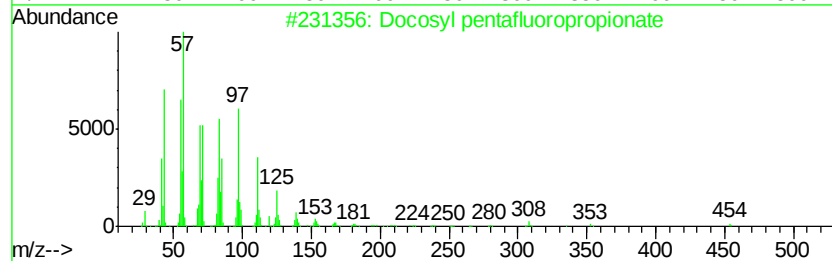
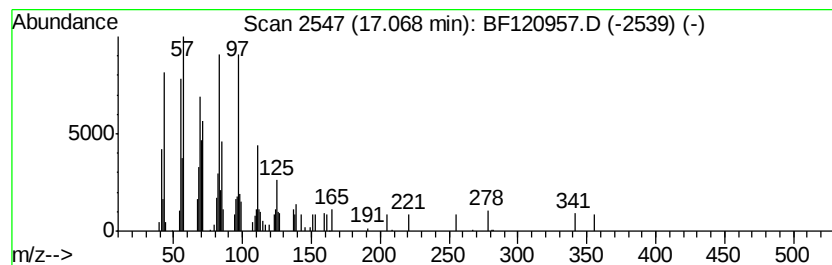
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Docosyl pentafluoropropionate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.07	2.58 ng	152775	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	90
2		Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	90
3		Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	90
4		Tetracosyl pentafluoropropionate	500	C27H49F5O2	1000351-81-3	90
5		Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072120\
Data File : BF120957.D
Acq On : 21 Jul 2020 16:46
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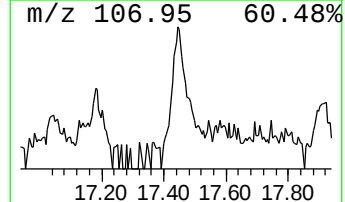
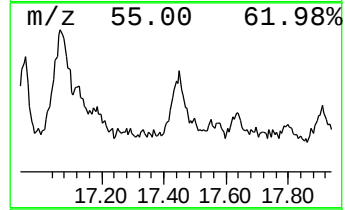
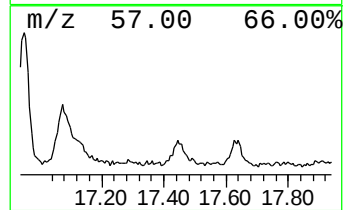
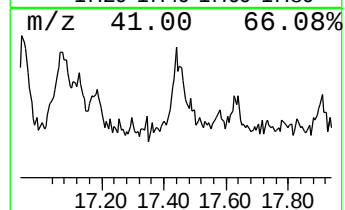
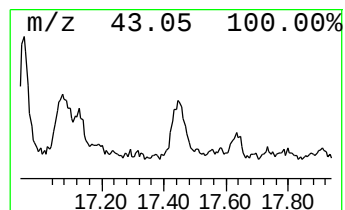
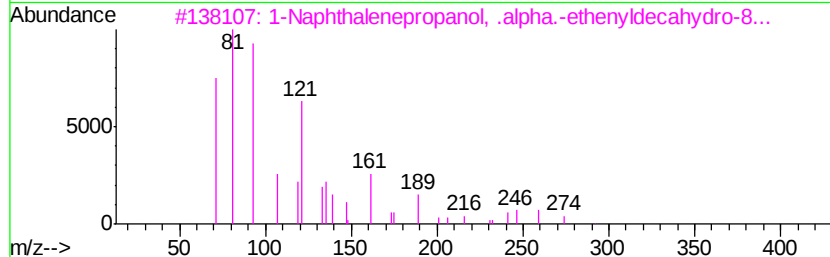
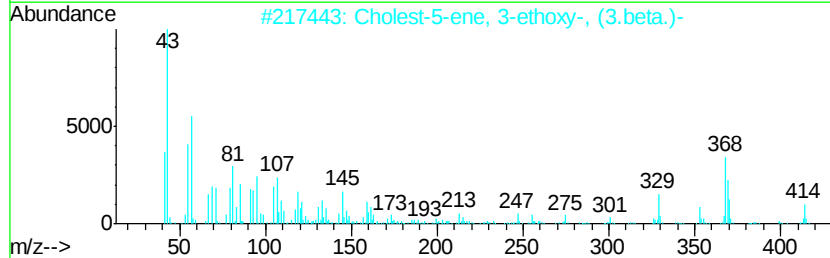
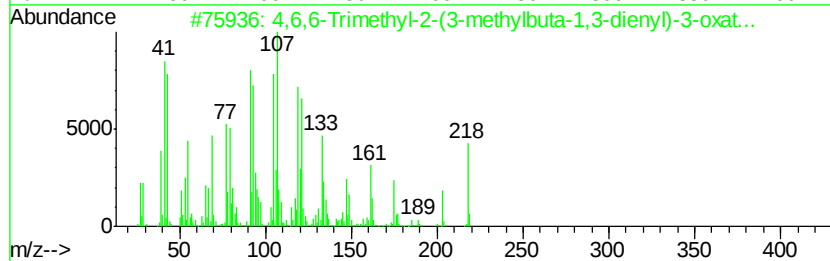
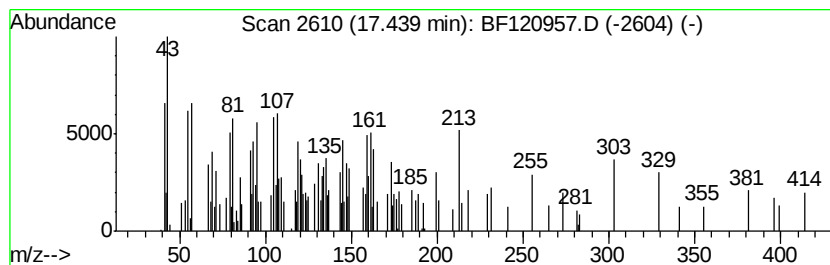
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 4,6,6-Trimethyl-2-(3-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.44	3.35 ng	198047	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4,6,6-Trimethyl-2-(3-methylbuta-...	218	C15H22O	1000190-22-2	55
2		Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	46
3		1-Naphthalenepropanol, .alpha.-e...	292	C19H32O2	055649-42-8	17
4		Acetic acid, 3-hydroxy-6-isoprop...	278	C17H26O3	1000185-44-8	16
5		1H-Cyclopenta[a]phenanthrene-16-...	398	C24H34N2O3	1000337-25-3	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072120\
 Data File : BF120957.D
 Acq On : 21 Jul 2020 16:46
 Operator : JU/CG
 Sample : L3354-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BACK-YARD

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
(+)-2-Bornanone	7.84	3.4	ng	218677	2	8.10	1304300	20.0
Safrrole	8.72	43.1	ng	2810300	2	8.10	1304300	20.0
2,4,4,6,6,8,8-Hep...	12.31	2.8	ng	231473	4	11.33	1648450	20.0
1-Heneicosanol	13.83	10.8	ng	813897	5	13.97	1514020	20.0
Cyclopentadecane	14.47	2.5	ng	192177	5	13.97	1514020	20.0
Octadecanal	14.89	2.7	ng	161969	6	15.42	1182150	20.0
Nonadecane	15.08	5.1	ng	299747	6	15.42	1182150	20.0
Behenic alcohol	15.12	6.4	ng	380001	6	15.42	1182150	20.0
Heneicosane	15.88	7.2	ng	424716	6	15.42	1182150	20.0
1-Heptacosanol	15.96	3.8	ng	221802	6	15.42	1182150	20.0
Docosyl pentaflu...	17.07	2.6	ng	152775	6	15.42	1182150	20.0
4,6,6-Trimethyl-2...	17.44	3.4	ng	198047	6	15.42	1182150	20.0