

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032720\
 Data File : BF119621.D
 Acq On : 28 Mar 2020 17:16
 Operator : CG/JU
 Sample : L2035-03
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-054-B

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-BF031820.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	2.128	3	7	12	rVB	349990	446620	5.11%	0.478%
2	2.234	21	25	31	rBV2	161749	203880	2.33%	0.218%
3	5.057	496	505	513	rBV	677603	909499	10.41%	0.973%
4	5.457	567	573	576	rBV	7833869	8733398	100.00%	9.344%
5	6.463	738	744	747	rBV	7014852	8621398	98.72%	9.224%
6	6.834	804	807	811	rBV	2579278	2164174	24.78%	2.315%
7	6.993	830	834	837	rBV	7098537	6479574	74.19%	6.933%
8	7.398	897	903	906	rBV	5222894	5457573	62.49%	5.839%
9	8.122	1021	1026	1028	rBV	3024929	2620098	30.00%	2.803%
10	8.698	1120	1124	1130	rBV2	170599	170794	1.96%	0.183%
11	9.198	1204	1209	1212	rBV	8633988	8613778	98.63%	9.216%
12	9.592	1272	1276	1280	rBV	513192	449214	5.14%	0.481%
13	9.734	1296	1300	1305	rVB	84958	108344	1.24%	0.116%
14	9.881	1318	1325	1328	rBV	3180418	2930566	33.56%	3.135%
15	9.910	1328	1330	1333	rVB	147315	115924	1.33%	0.124%
16	10.081	1355	1359	1363	rBV2	123268	121690	1.39%	0.130%
17	10.422	1414	1417	1424	rBV	106079	109092	1.25%	0.117%
18	10.669	1454	1459	1462	rBV	5061795	4692438	53.73%	5.021%
19	10.786	1476	1479	1484	rVB3	114823	144723	1.66%	0.155%
20	11.233	1549	1555	1558	rBV4	89508	127121	1.46%	0.136%
21	11.263	1558	1560	1565	rVB2	89820	88398	1.01%	0.095%
22	11.363	1574	1577	1580	rBV	2695354	2484347	28.45%	2.658%
23	11.386	1580	1581	1585	rVV	909836	609673	6.98%	0.652%
24	11.439	1585	1590	1593	rVB	233970	228399	2.62%	0.244%
25	11.592	1611	1616	1619	rBV	98870	121068	1.39%	0.130%
26	11.763	1642	1645	1651	rVB2	76611	92356	1.06%	0.099%
27	11.869	1659	1663	1665	rBV	137409	152137	1.74%	0.163%
28	11.898	1665	1668	1674	rVV	2419867	2191859	25.10%	2.345%
29	11.980	1678	1682	1684	rBV2	195131	191642	2.19%	0.205%
30	12.069	1694	1697	1700	rVB2	105178	91661	1.05%	0.098%
31	12.416	1753	1756	1759	rBV	109063	106318	1.22%	0.114%
32	12.451	1759	1762	1763	rVV2	222400	208671	2.39%	0.223%
33	12.469	1763	1765	1768	rVB3	382243	322968	3.70%	0.346%
34	12.580	1780	1784	1786	rBV	1325852	1195610	13.69%	1.279%

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35	12.604	1786	1788	1796	rVB5	175777	262163	3.00%	0.280%
36	12.686	1798	1802	1808	rBV	1220589	1372337	15.71%	1.468%
37	12.780	1815	1818	1819	rBV	337486	320373	3.67%	0.343%
38	12.810	1820	1823	1825	rVB	1061993	955535	10.94%	1.022%
39	12.957	1844	1848	1851	rBV	6446090	5865163	67.16%	6.275%
40	13.133	1876	1878	1881	rVB	192961	186789	2.14%	0.200%
41	13.204	1885	1890	1892	rBV3	189777	260670	2.98%	0.279%
42	13.239	1893	1896	1898	rBV2	138719	122657	1.40%	0.131%
43	13.433	1927	1929	1934	rBV5	126858	136699	1.57%	0.146%
44	13.486	1935	1938	1941	rVB	207563	168609	1.93%	0.180%
45	13.533	1943	1946	1949	rVB	167502	138182	1.58%	0.148%
46	13.863	1999	2002	2011	rVB3	922635	1260228	14.43%	1.348%
47	14.010	2022	2027	2029	rBV2	2152337	2617301	29.97%	2.800%
48	14.033	2029	2031	2037	rVB	677929	535807	6.14%	0.573%
49	14.180	2052	2056	2061	rBV	1158924	1116702	12.79%	1.195%
50	14.486	2104	2108	2111	rBV	1981148	1848887	21.17%	1.978%
51	14.516	2111	2113	2116	rBV4	171059	212592	2.43%	0.227%
52	14.792	2156	2160	2164	rBV	2534593	2436758	27.90%	2.607%
53	15.045	2200	2203	2206	rBV	564448	751657	8.61%	0.804%
54	15.133	2214	2218	2224	rBV	2528250	2886093	33.05%	3.088%
55	15.345	2252	2254	2261	rVB2	351828	480974	5.51%	0.515%
56	15.410	2262	2265	2269	rVV	444630	544951	6.24%	0.583%
57	15.474	2272	2276	2279	rVV	1313610	1594847	18.26%	1.706%
58	15.515	2279	2283	2290	rVB	2012799	2735037	31.32%	2.926%
59	15.957	2353	2358	2363	rBV	1378012	2088199	23.91%	2.234%
60	16.462	2440	2444	2456	rVB2	634569	1260568	14.43%	1.349%

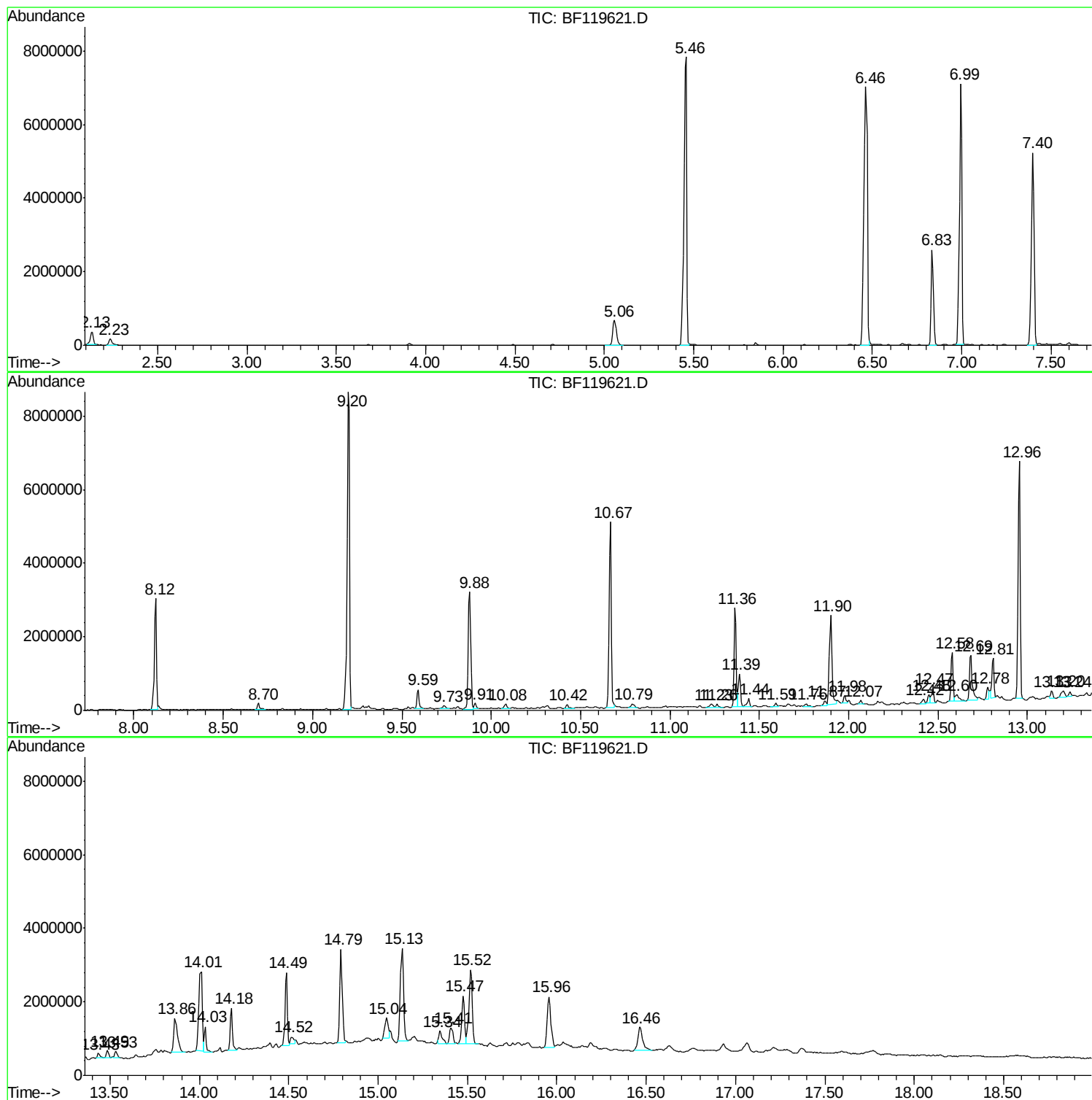
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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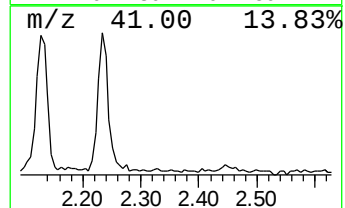
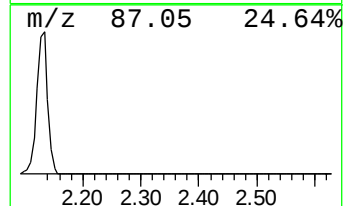
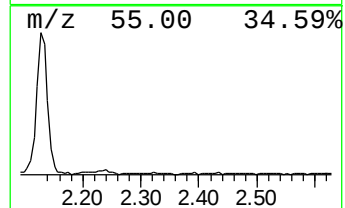
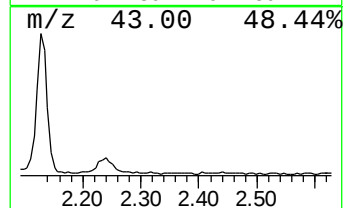
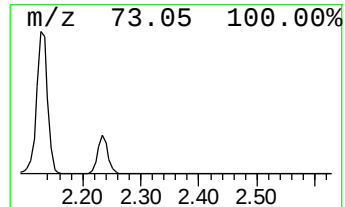
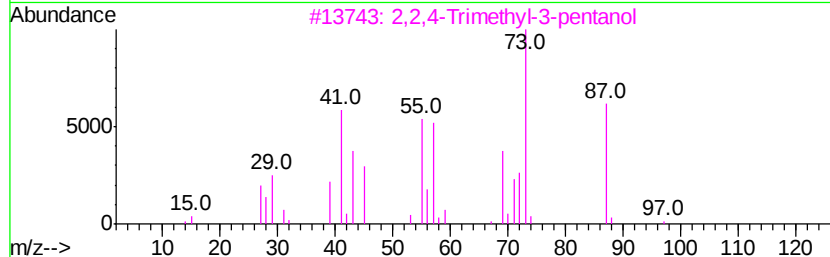
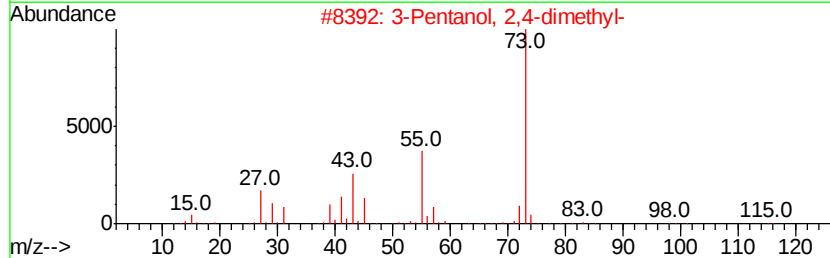
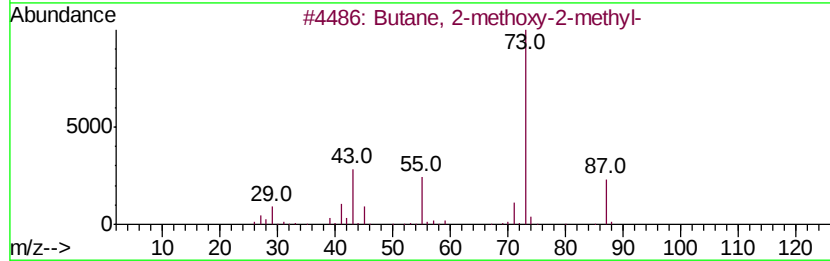
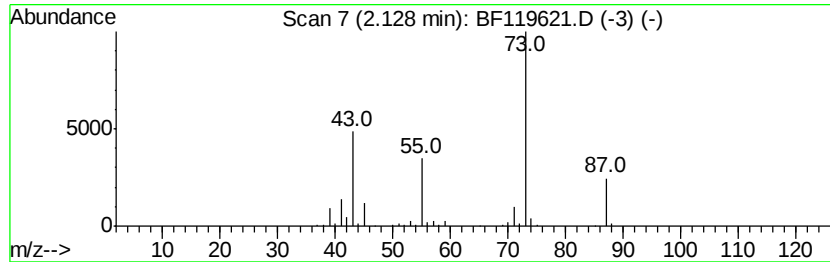
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	4.13 ng	446620	1,4-Dichlorobenzene-d4	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	38
5		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	17



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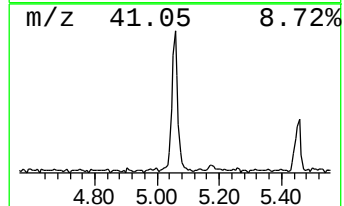
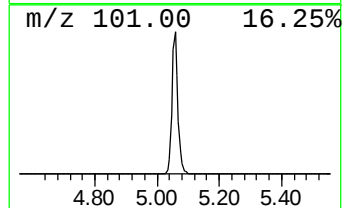
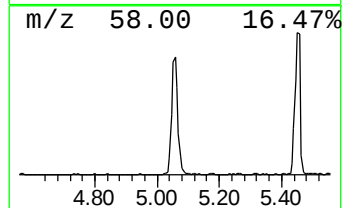
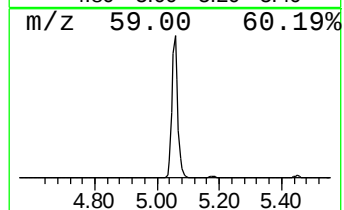
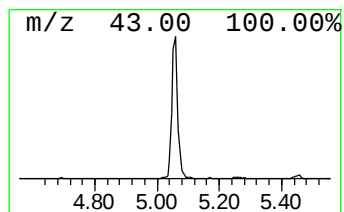
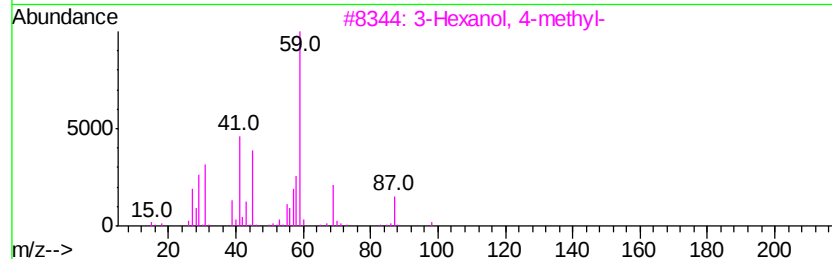
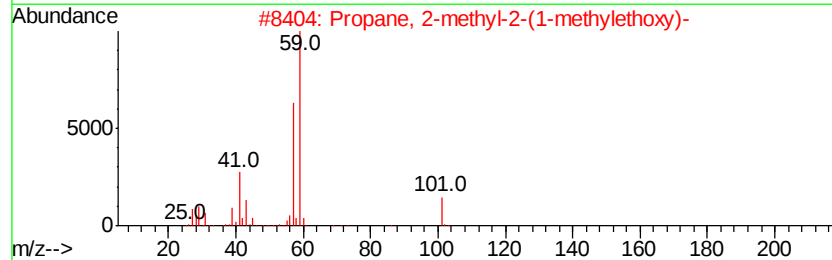
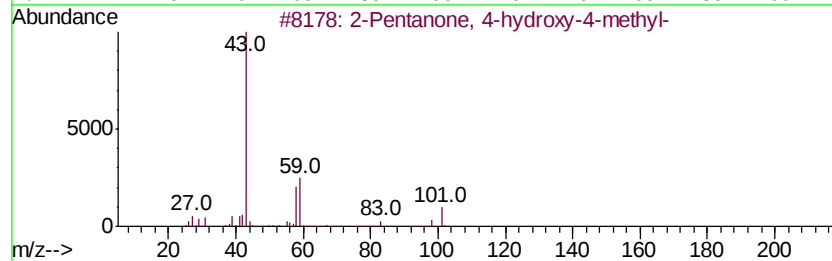
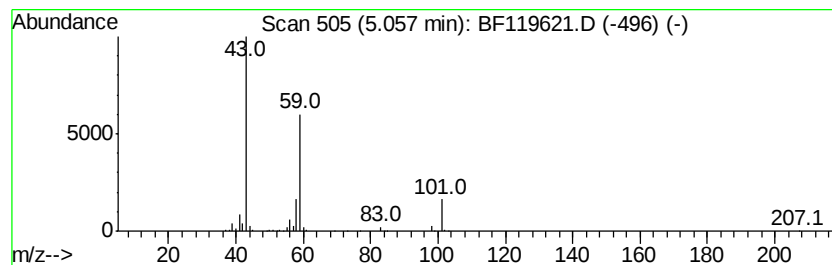
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	8.41 ng	909499	1,4-Dichlorobenzene-d4	6.83
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116 C6H12O2	000123-42-2	59
2	Propane, 2-methyl-2-(1-methyleth...	116 C7H16O	017348-59-3	42
3	3-Hexanol, 4-methyl-	116 C7H16O	000615-29-2	33
4	2-Hexanol, 2-methyl-	116 C7H16O	000625-23-0	33
5	Acetic acid, cyano-, 1,1-dimethy...	141 C7H11NO2	001116-98-9	25



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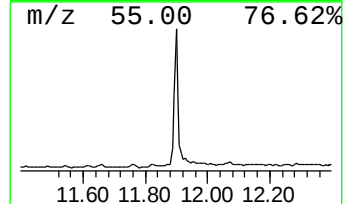
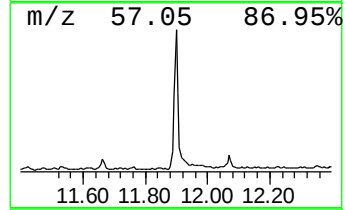
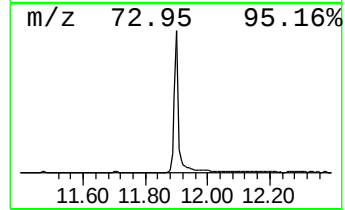
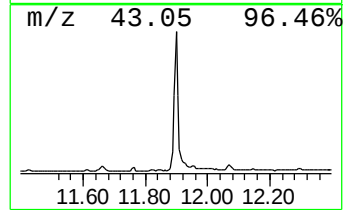
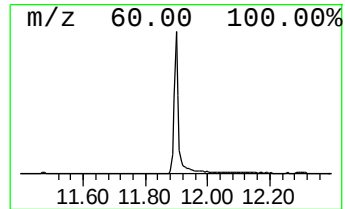
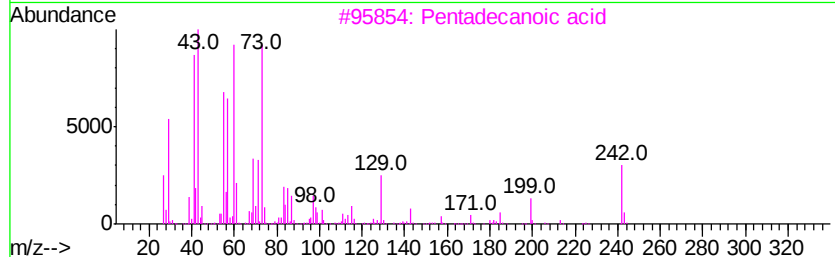
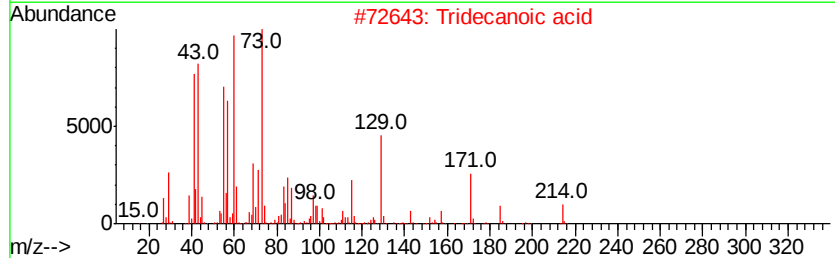
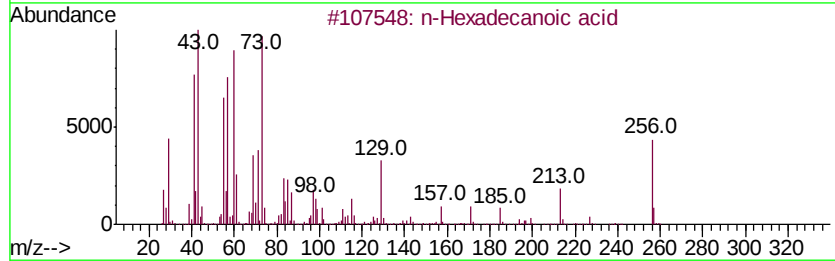
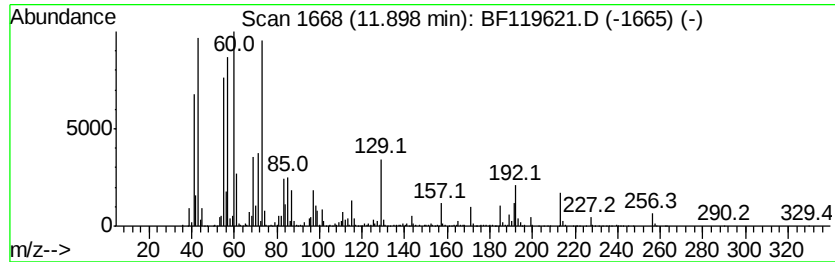
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	17.65 ng	2191860	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Tridecanoic acid	214	C13H26O2	000638-53-9	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Undecanoic acid	186	C11H22O2	000112-37-8	64



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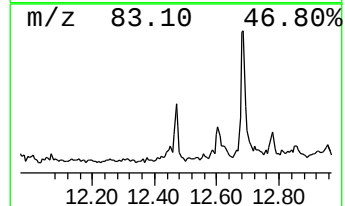
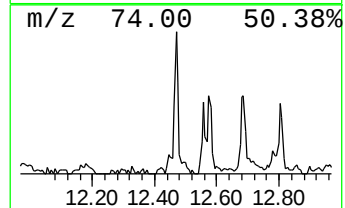
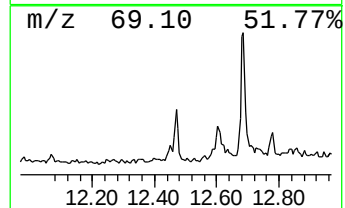
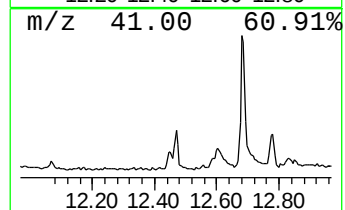
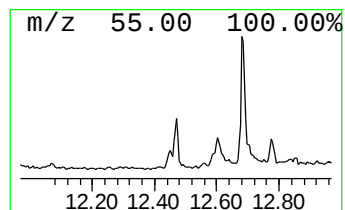
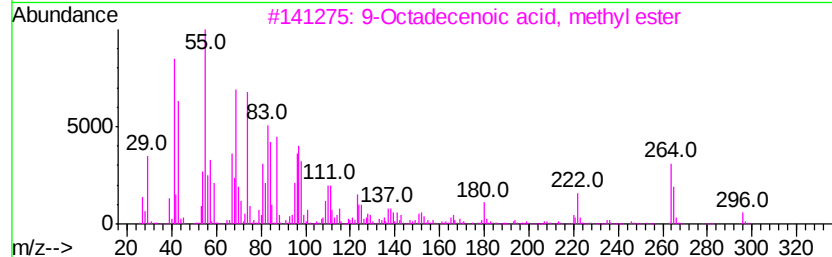
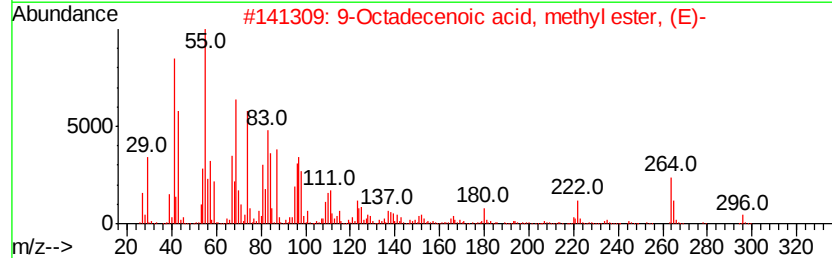
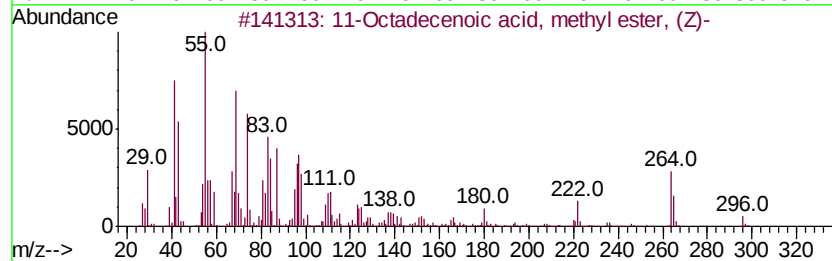
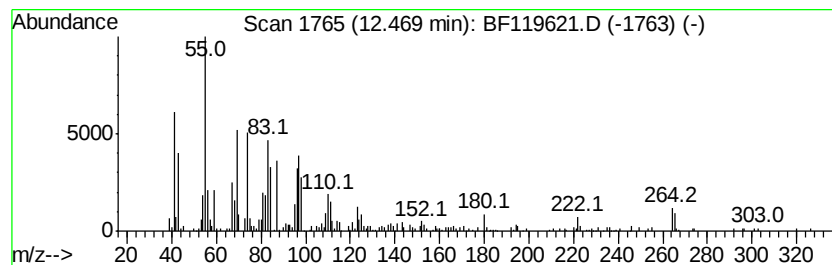
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 11-Octadecenoic acid, methyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	2.60 ng	322968	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-63-9	99
2		9-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-62-8	99
3		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
4		trans-13-Octadecenoic acid, methyl ester...	296	C19H36O2	1000333-61-3	99
5		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	98



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FK-GF-02-054-B

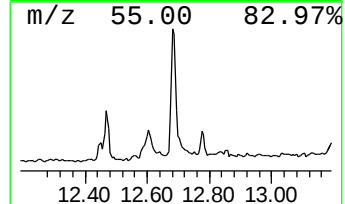
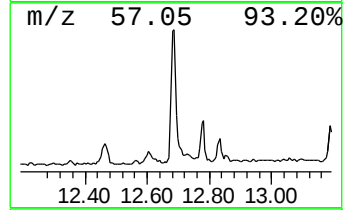
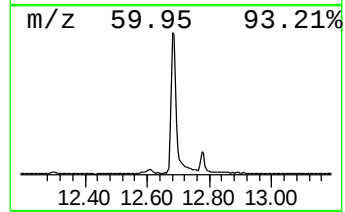
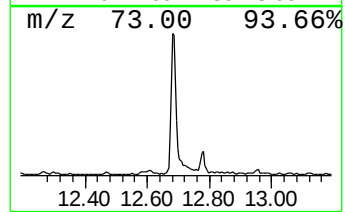
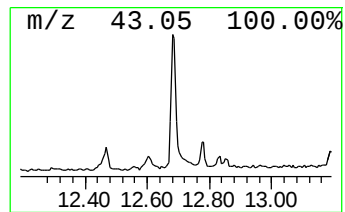
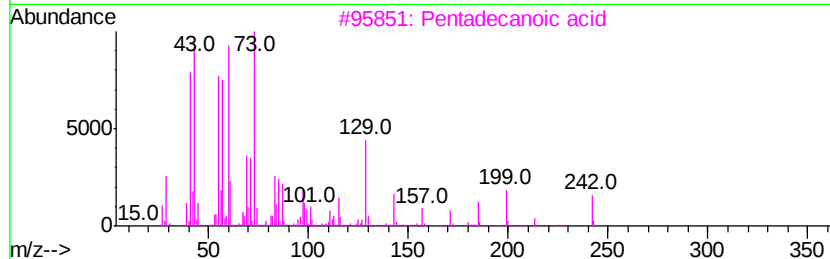
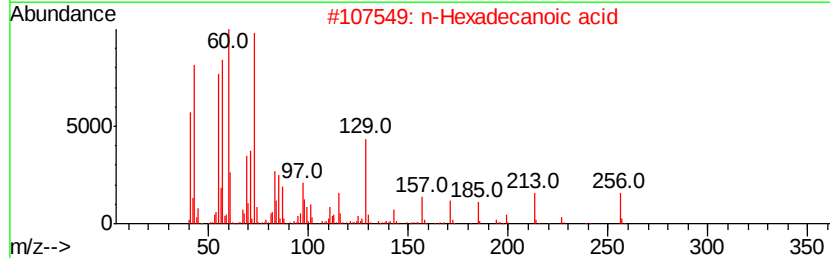
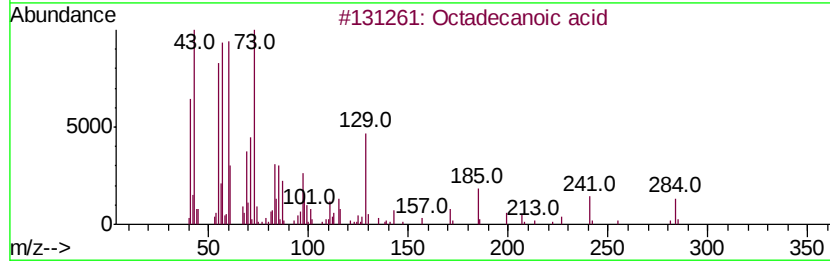
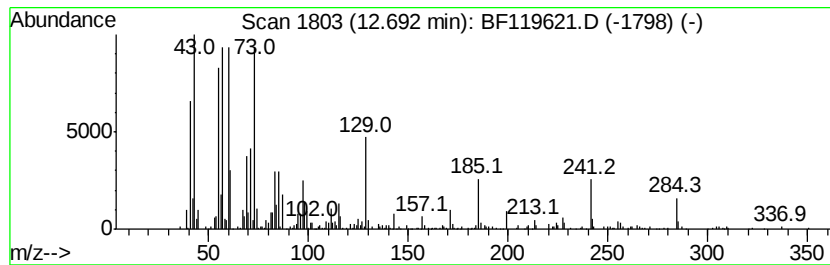
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	10.49 ng	1372340	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		n-Decanoic acid	172	C10H20O2	000334-48-5	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
Operator : CG/JU
Sample : L2035-03
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-054-B

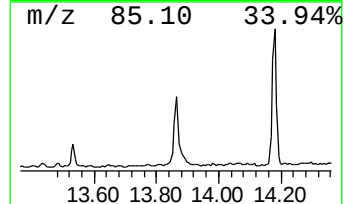
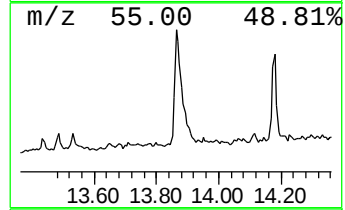
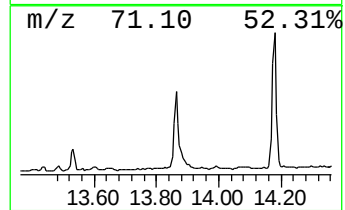
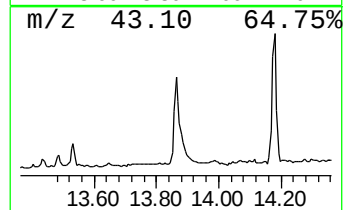
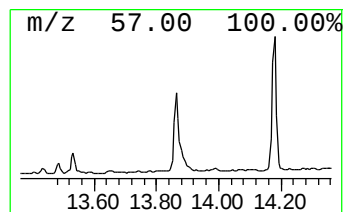
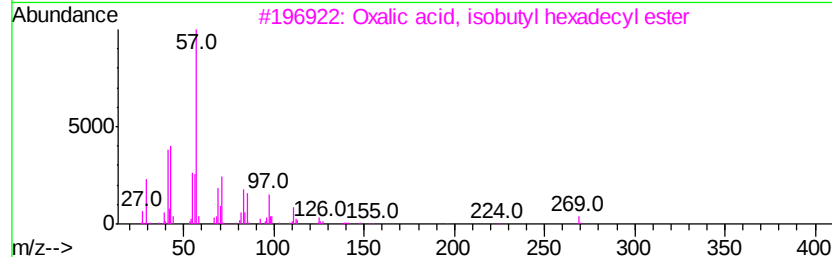
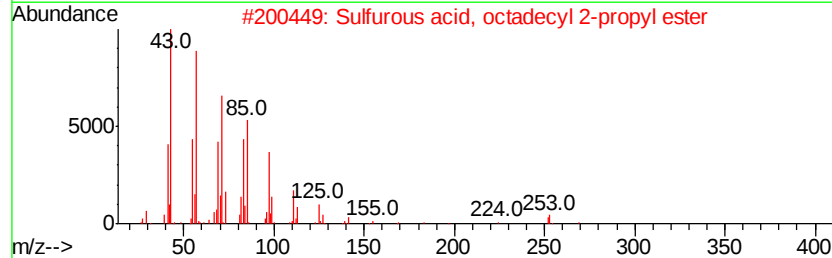
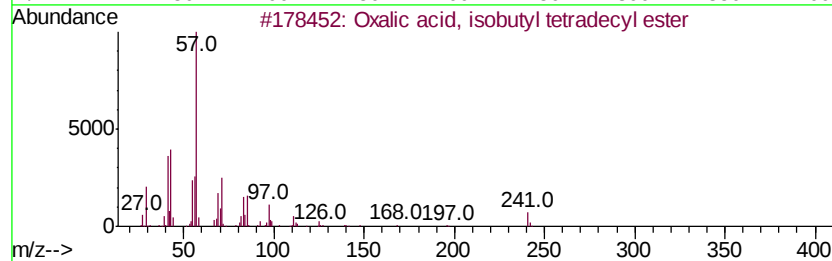
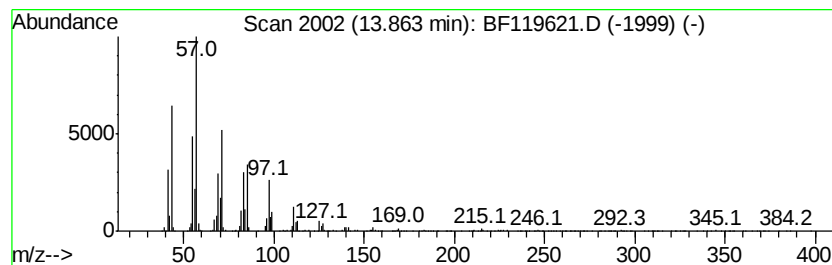
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxalic acid, isobutyl tetra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	9.63 ng	1260230	Chrysene-d12	14.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	87
2			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	87
3			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	76
4			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	76
5			Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
Operator : CG/JU
Sample : L2035-03
Misc :
ALS Vial : 51 Sample Multiplier: 1

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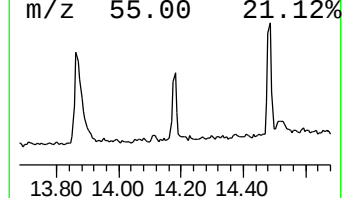
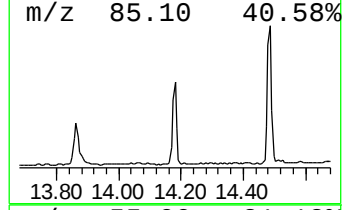
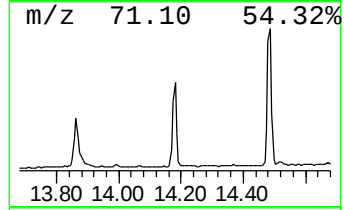
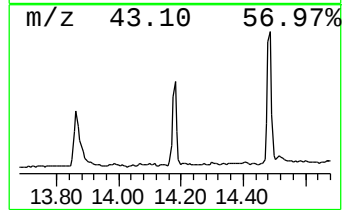
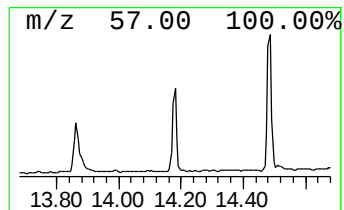
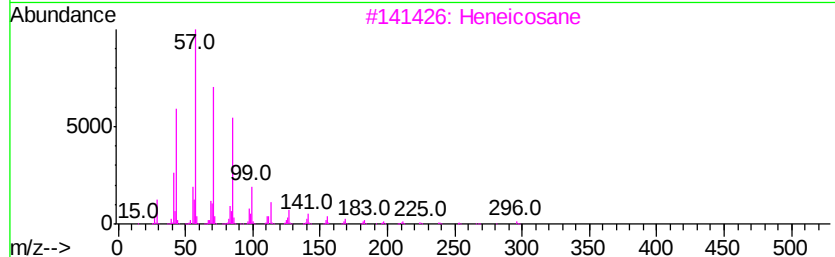
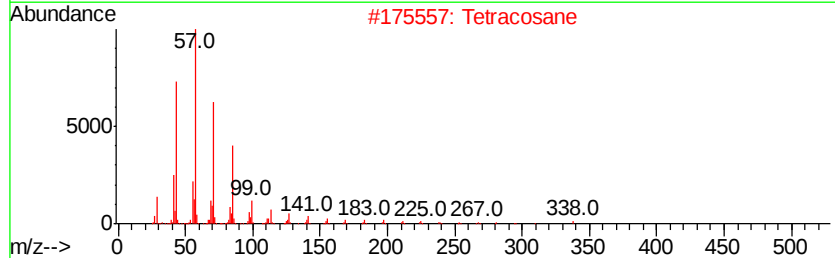
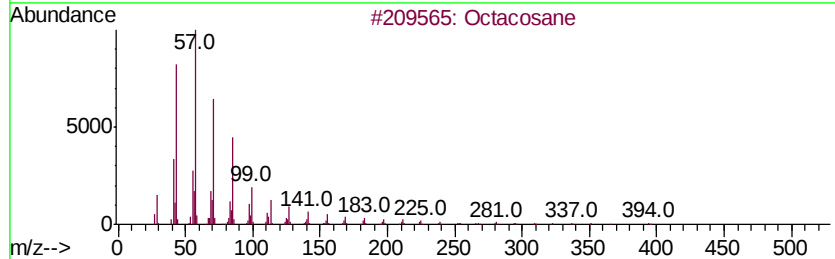
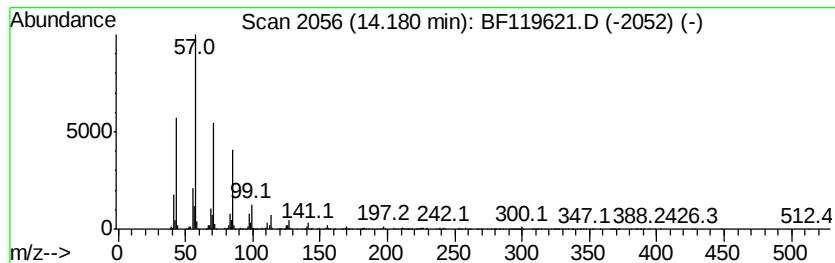
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	8.53 ng	1116700	Chrysene-d12	14.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	93
2			Tetracosane	338	C24H50	000646-31-1	91
3			Heneicosane	296	C21H44	000629-94-7	90
4			Octadecane	254	C18H38	000593-45-3	90
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
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Sample : L2035-03
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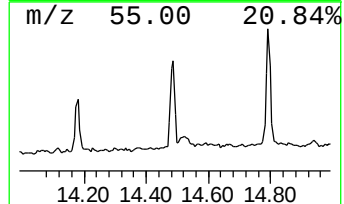
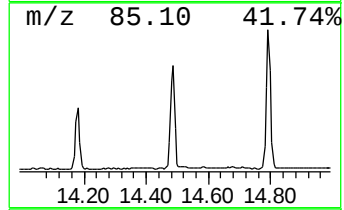
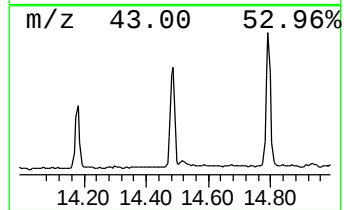
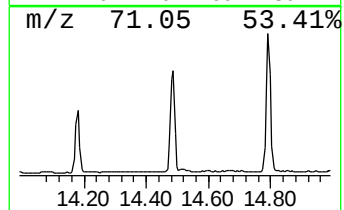
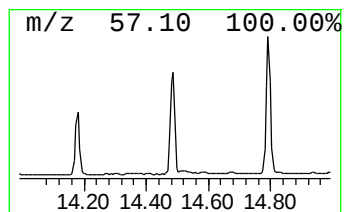
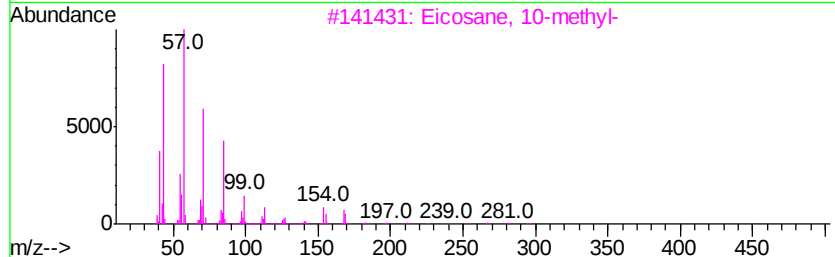
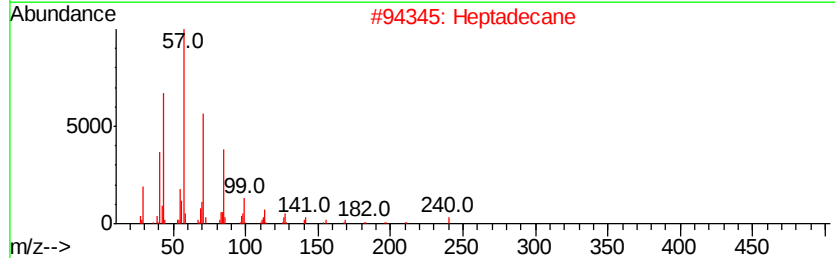
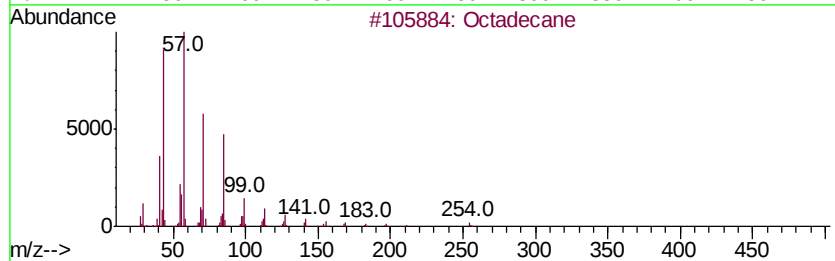
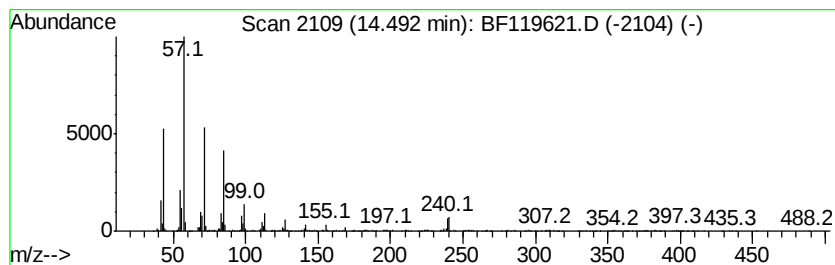
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.49	14.13 ng	1848890	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	95
2	Heptadecane	240 C17H36	000629-78-7	93
3	Eicosane, 10-methyl-	296 C21H44	054833-23-7	93
4	1-Iodo-2-methylundecane	296 C12H25I	073105-67-6	91
5	Tetracosane	338 C24H50	000646-31-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
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Instrument :
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ClientSampleId :
FK-GF-02-054-B

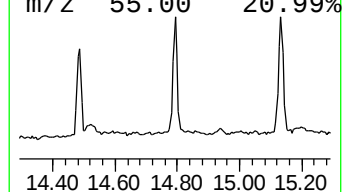
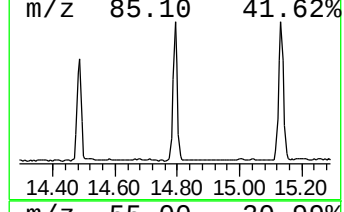
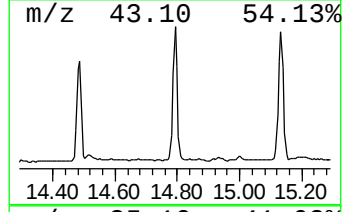
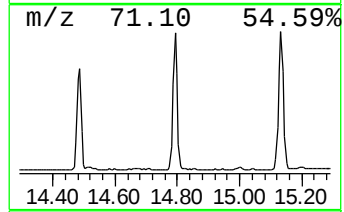
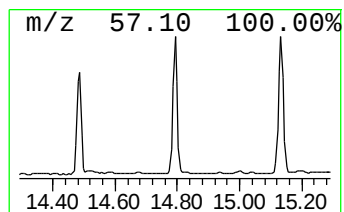
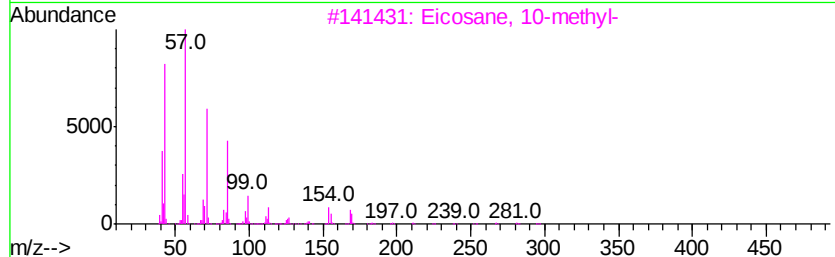
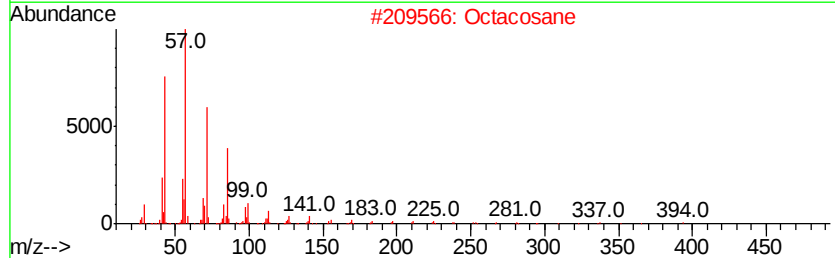
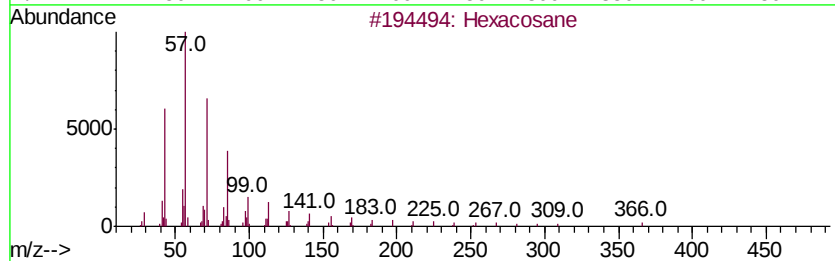
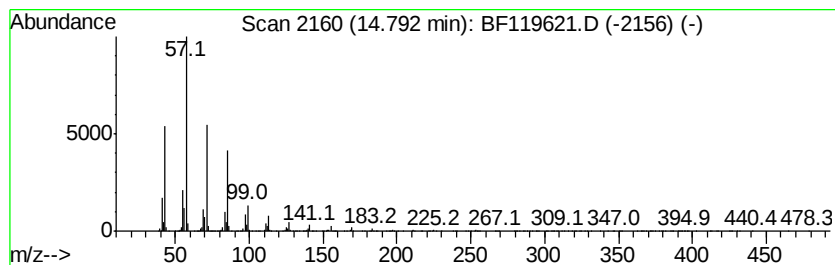
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	30.56 ng	2436760	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	95
2		Octacosane	394	C28H58	000630-02-4	95
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentacosane	352	C25H52	000629-99-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
Operator : CG/JU
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Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FK-GF-02-054-B

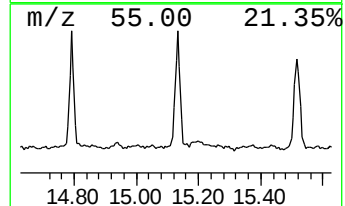
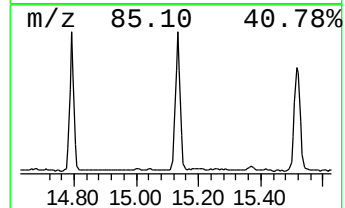
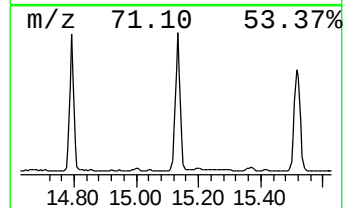
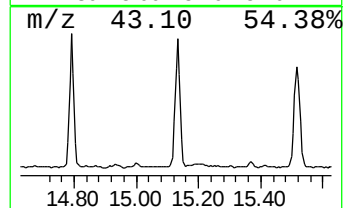
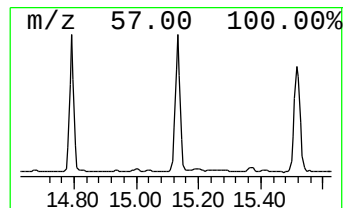
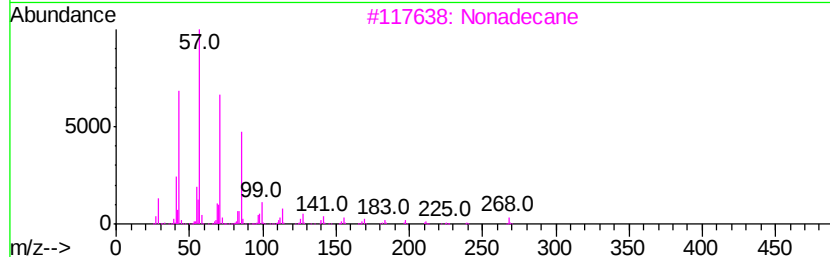
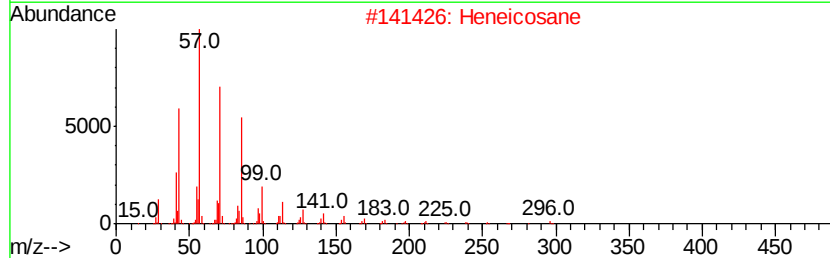
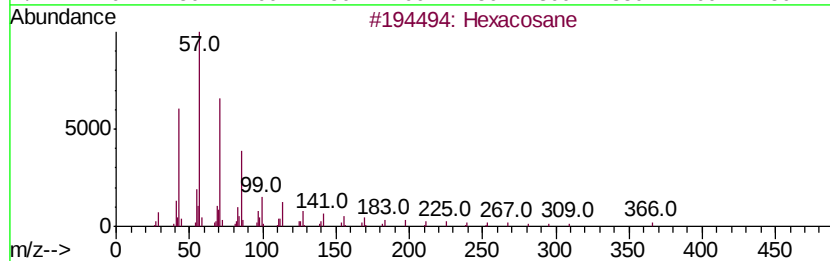
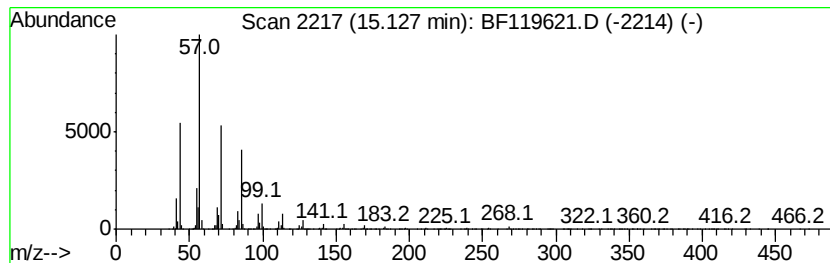
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	36.19 ng	2886090	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	92
2		Heneicosane	296	C21H44	000629-94-7	90
3		Nonadecane	268	C19H40	000629-92-5	90
4		Triacontane	422	C30H62	000638-68-6	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
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ClientSampleId :
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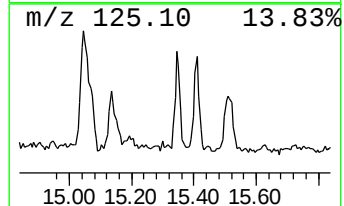
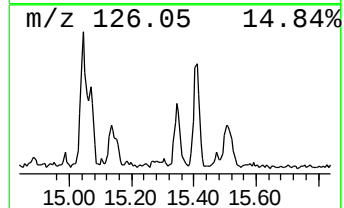
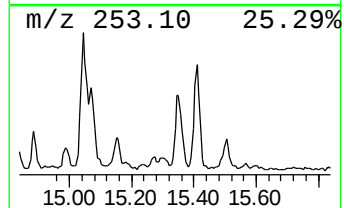
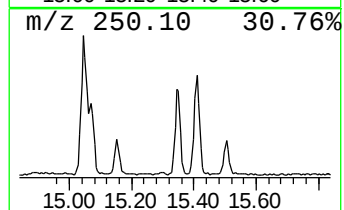
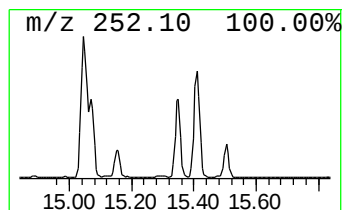
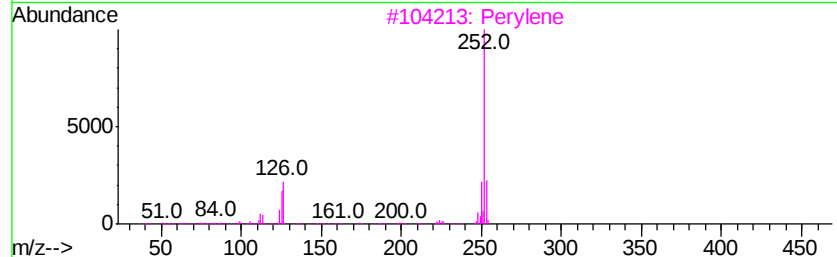
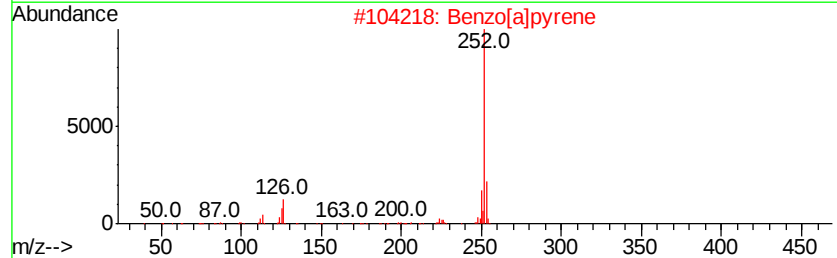
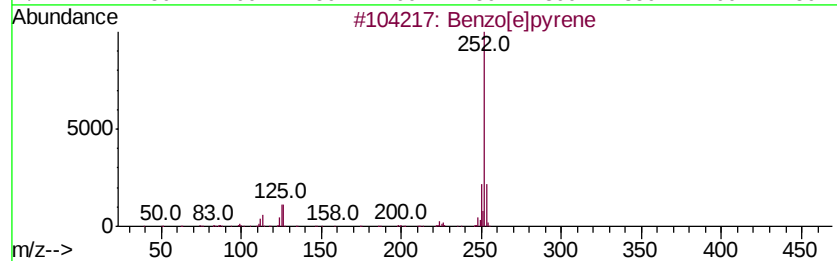
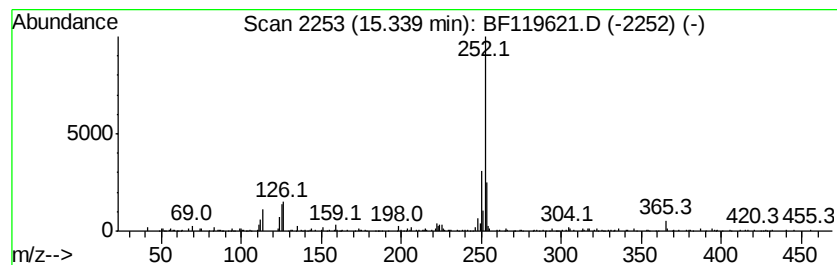
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	6.03 ng	480974	Perylene-d12	15.47

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
Operator : CG/JU
Sample : L2035-03
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-054-B

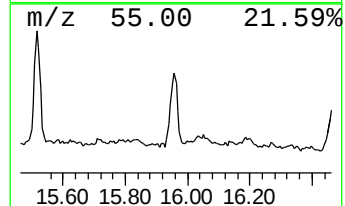
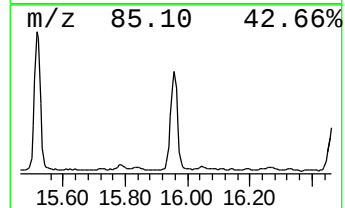
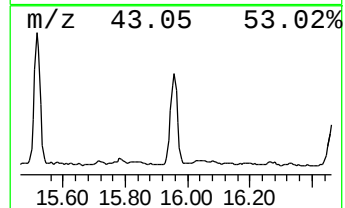
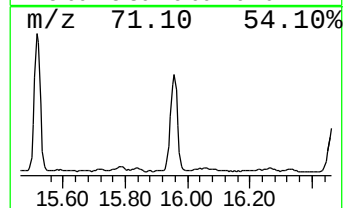
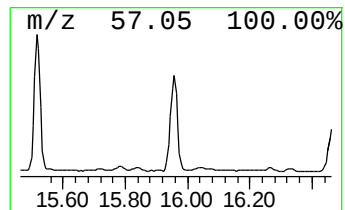
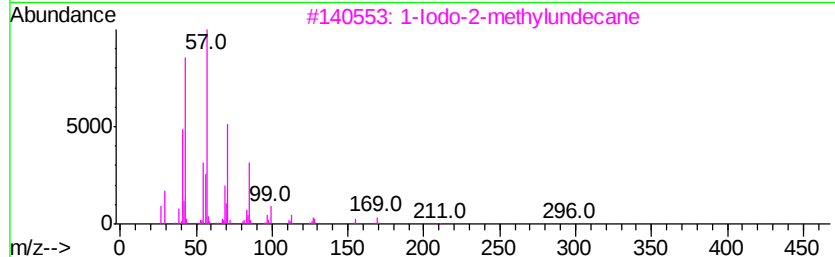
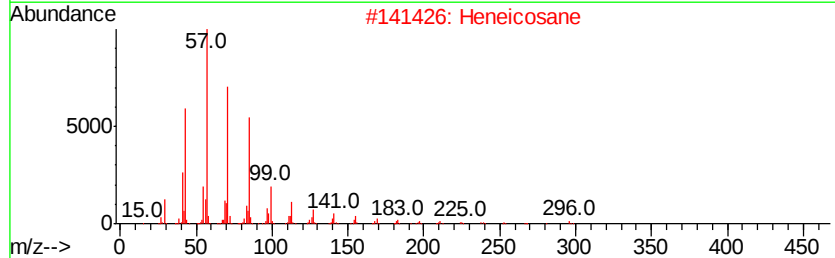
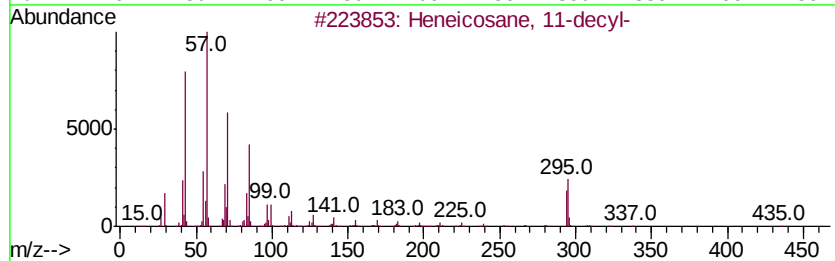
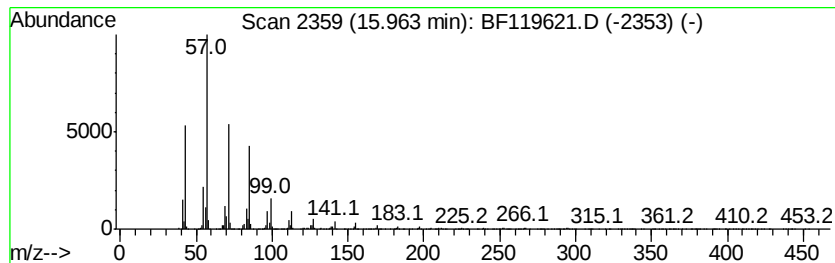
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane, 11-decyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	26.19 ng	2088200	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4		Eicosane	282	C20H42	000112-95-8	83
5		Hentriacontane	437	C31H64	000630-04-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
Data File : BF119621.D
Acq On : 28 Mar 2020 17:16
Operator : CG/JU
Sample : L2035-03
Misc :
ALS Vial : 51 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-02-054-B

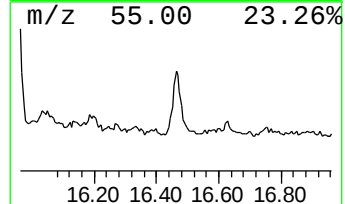
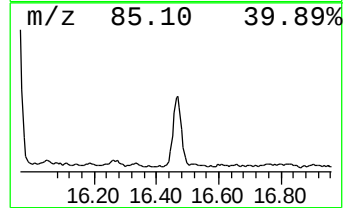
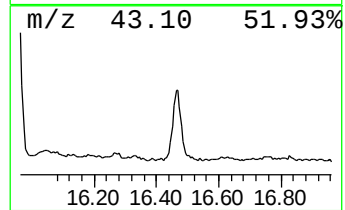
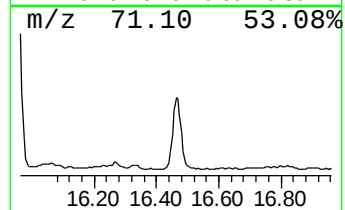
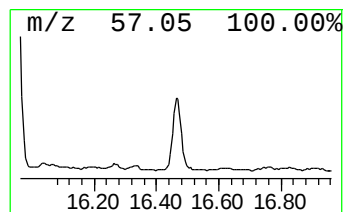
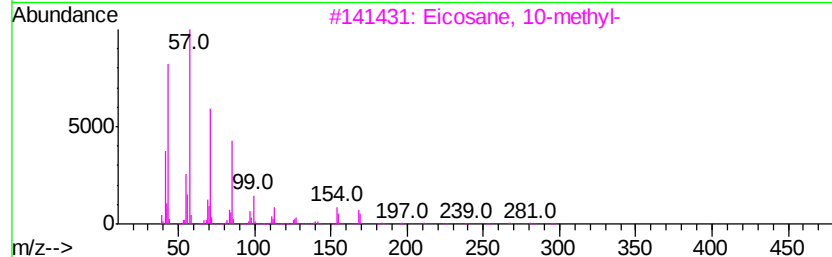
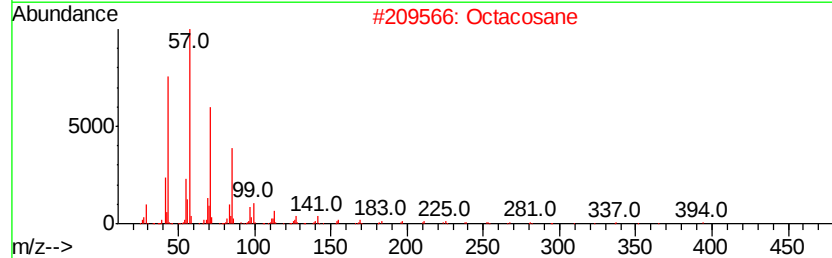
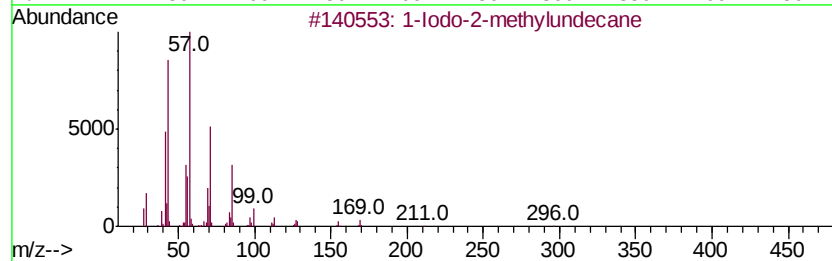
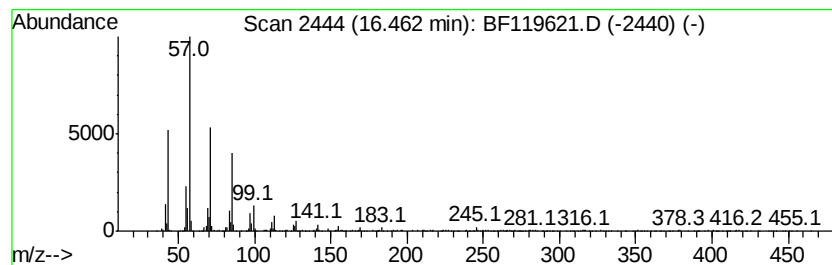
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Iodo-2-methylundecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	15.81 ng	1260570	Perylene-d12	15.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
2		Octacosane	394	C28H58	000630-02-4	90
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
4		Pentacosane	352	C25H52	000629-99-2	87
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032720\
 Data File : BF119621.D
 Acq On : 28 Mar 2020 17:16
 Operator : CG/JU
 Sample : L2035-03
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-054-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF031820.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
Butane, 2-methoxy...	2.13	4.1 ng		446620	1	6.83	2164170	20.0
2-Pentanone, 4-hy...	5.06	8.4 ng		909499	1	6.83	2164170	20.0
n-Hexadecanoic acid	11.90	17.6 ng		2191860	4	11.36	2484350	20.0
11-Octadecenoic a...	12.47	2.6 ng		322968	4	11.36	2484350	20.0
Octadecanoic acid	12.69	10.5 ng		1372340	5	14.01	2617300	20.0
Oxalic acid, isob...	13.86	9.6 ng		1260230	5	14.01	2617300	20.0
Octacosane	14.18	8.5 ng		1116700	5	14.01	2617300	20.0
Octadecane	14.49	14.1 ng		1848890	5	14.01	2617300	20.0
Hexacosane	14.79	30.6 ng		2436760	6	15.47	1594850	20.0
Heneicosane	15.13	36.2 ng		2886090	6	15.47	1594850	20.0
Benzo[e]pyrene	15.34	6.0 ng		480974	6	15.47	1594850	20.0
Heneicosane, 11-d...	15.96	26.2 ng		2088200	6	15.47	1594850	20.0
1-Iodo-2-methylun...	16.46	15.8 ng		1260570	6	15.47	1594850	20.0