

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114128.D
 Acq On : 10 May 2019 18:48
 Operator : JU/SJ
 Sample : K2762-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P021-SS002-0206-01

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_F\METHODS\8270-BF051019.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	2.152	5	11	19	rVB	1281687	1713446	23.01%	2.693%
2	5.481	572	577	580	rBV	7212433	6912202	92.82%	10.865%
3	6.487	743	748	751	rBV	7304843	7446655	100.00%	11.705%
4	6.622	767	771	775	rBV	7091453	7151953	96.04%	11.242%
5	6.851	806	810	814	rBV	2358094	1908042	25.62%	2.999%
6	7.010	833	837	840	rBV	5195423	4366262	58.63%	6.863%
7	7.410	900	905	908	rBV	4036493	3639815	48.88%	5.721%
8	8.133	1023	1028	1031	rBV	2327673	2069066	27.79%	3.252%
9	8.575	1100	1103	1117	rVB	162884	202267	2.72%	0.318%
10	9.204	1205	1210	1214	rBV	6637196	6150148	82.59%	9.667%
11	9.598	1273	1277	1280	rBV	992246	825998	11.09%	1.298%
12	9.886	1319	1326	1329	rBV	2762732	2411537	32.38%	3.791%
13	10.669	1455	1459	1463	rBV	3930714	3774724	50.69%	5.933%
14	11.369	1574	1578	1581	rBV	2592412	2245616	30.16%	3.530%
15	11.892	1664	1667	1679	rBV	533369	606194	8.14%	0.953%
16	12.680	1797	1801	1804	rBV	131000	158917	2.13%	0.250%
17	12.951	1842	1847	1850	rBV	5029557	4390356	58.96%	6.901%
18	13.851	1996	2000	2012	rBV2	317595	496096	6.66%	0.780%
19	14.010	2022	2027	2030	rBV	1932157	1759059	23.62%	2.765%
20	14.180	2052	2056	2064	rVB	147166	166457	2.24%	0.262%
21	14.498	2106	2110	2114	rBV	227330	237076	3.18%	0.373%
22	14.804	2159	2162	2166	rVB	321349	293380	3.94%	0.461%
23	14.880	2172	2175	2179	rVB	250731	231272	3.11%	0.364%
24	15.145	2216	2220	2224	rBV	413166	464610	6.24%	0.730%
25	15.492	2274	2279	2282	rBV	1390264	1727031	23.19%	2.715%
26	15.521	2282	2284	2290	rVB	446985	510907	6.86%	0.803%
27	15.957	2353	2358	2364	rBV	432433	648645	8.71%	1.020%
28	16.462	2439	2444	2450	rVB	226193	388994	5.22%	0.611%
29	17.051	2541	2544	2552	rVB3	101905	194147	2.61%	0.305%
30	18.045	2706	2713	2720	rVB7	231808	527449	7.08%	0.829%

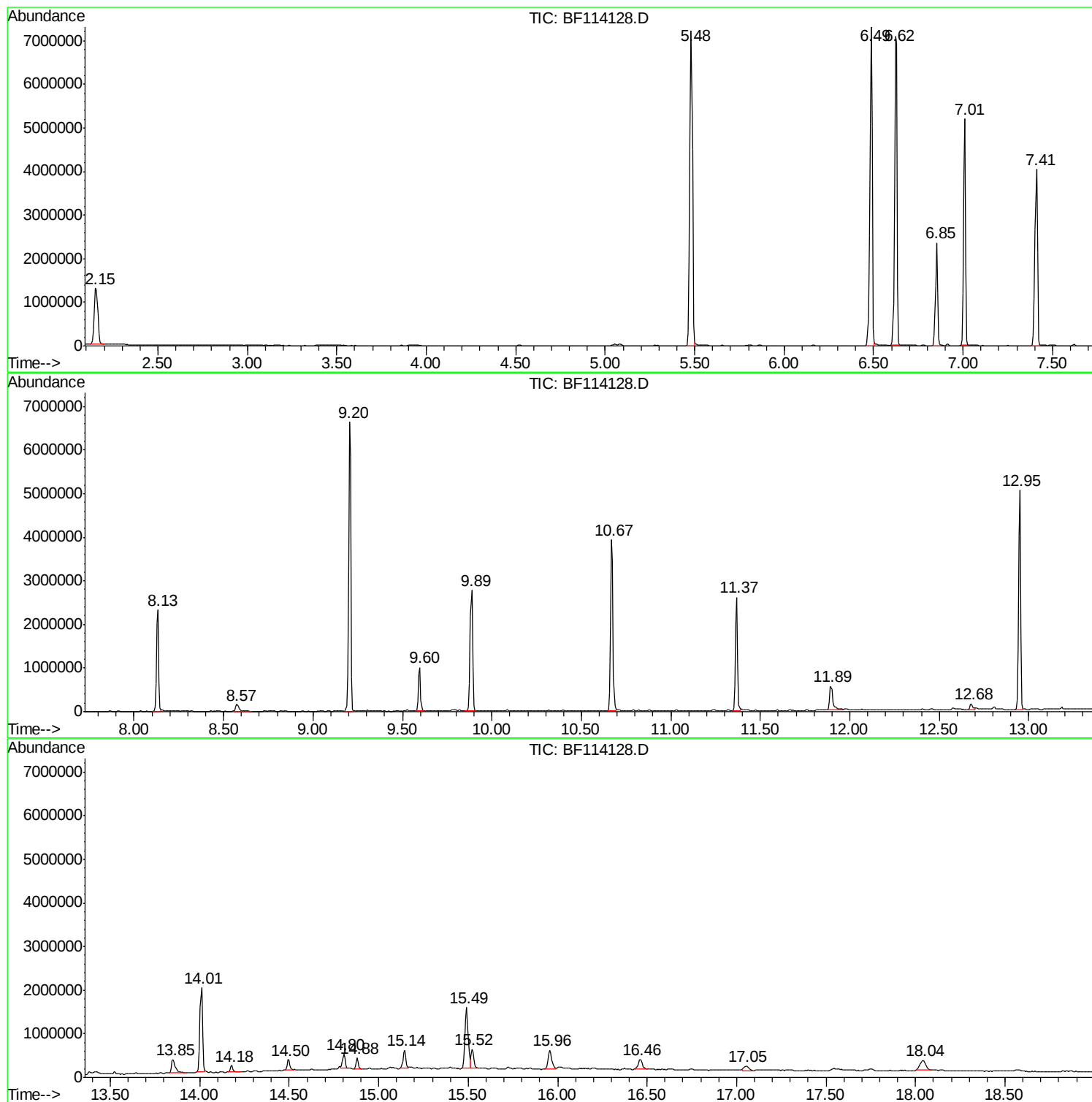
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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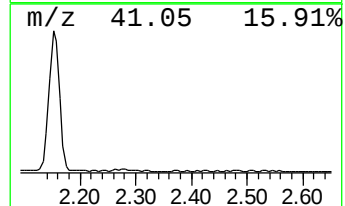
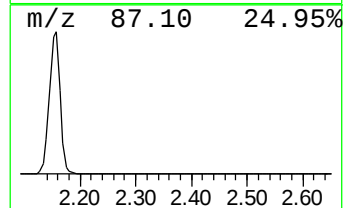
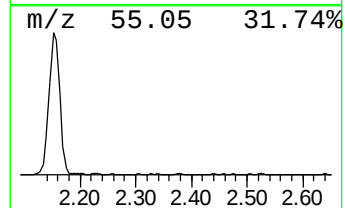
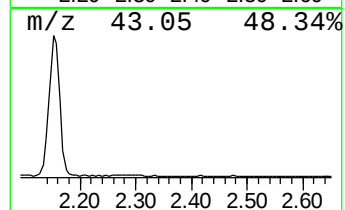
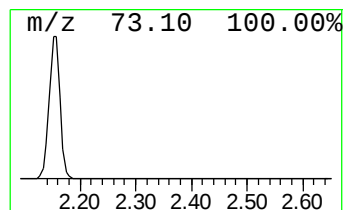
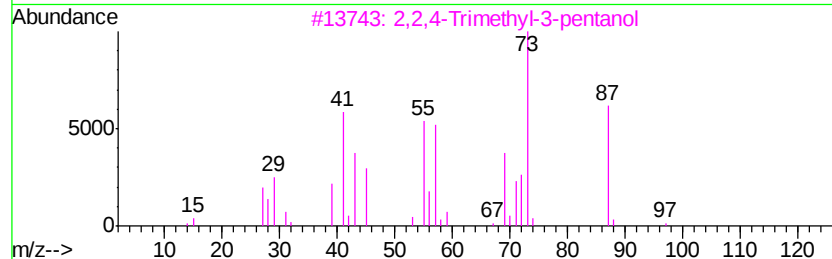
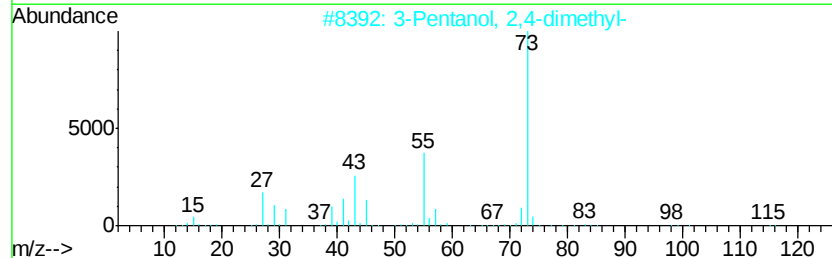
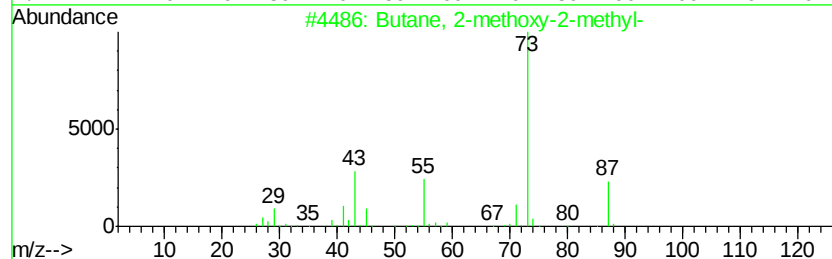
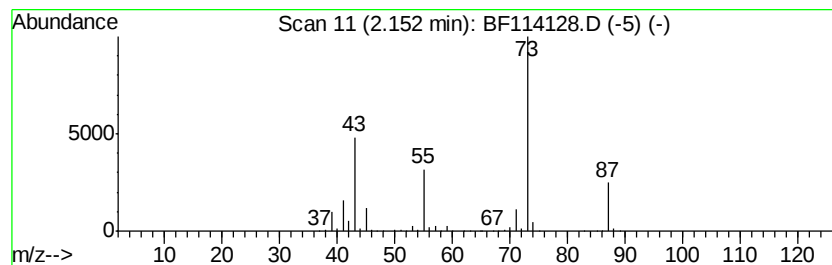
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	17.96 ng	1713450	1,4-Dichlorobenzene-d4	6.85

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	40
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23



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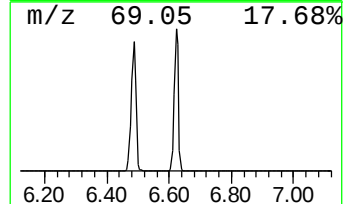
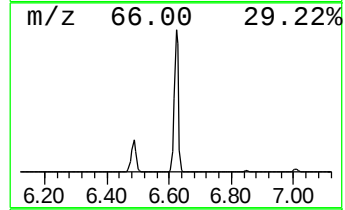
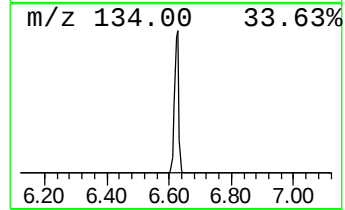
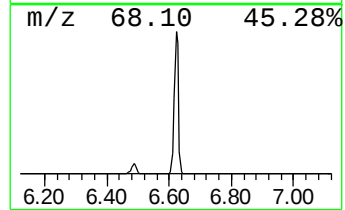
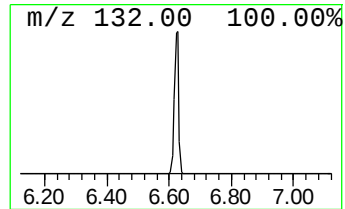
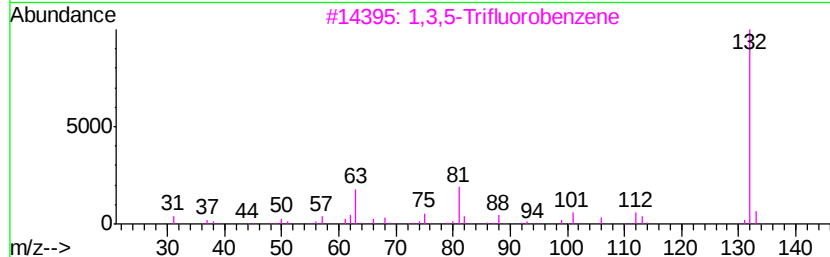
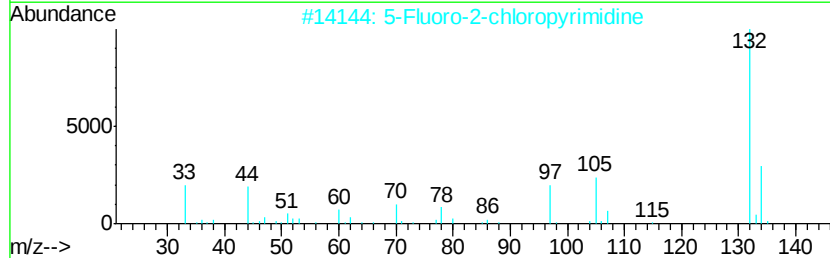
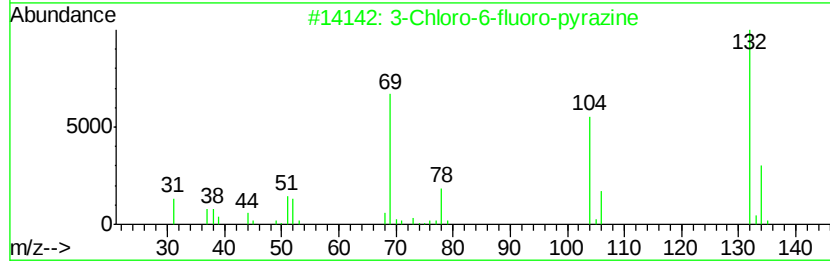
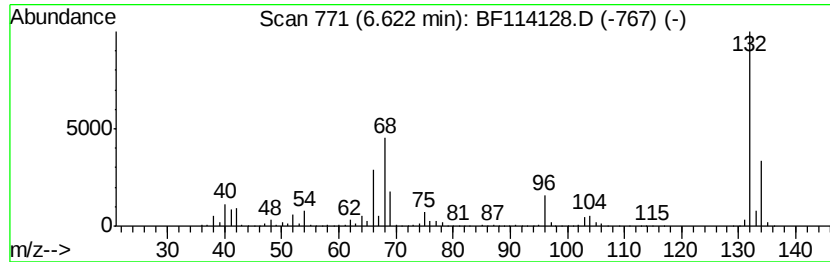
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	74.97 ng	7151950	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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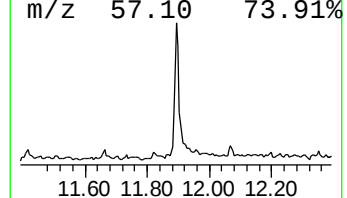
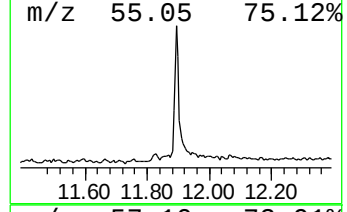
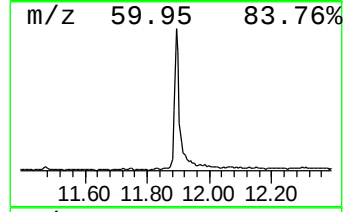
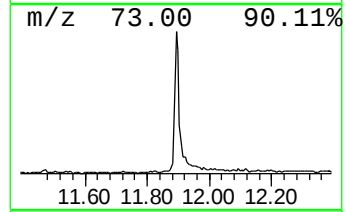
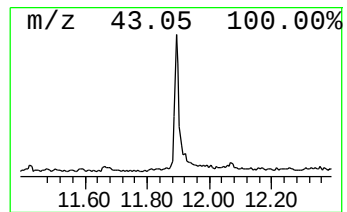
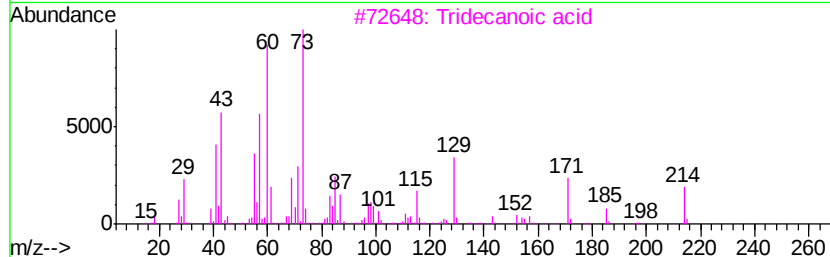
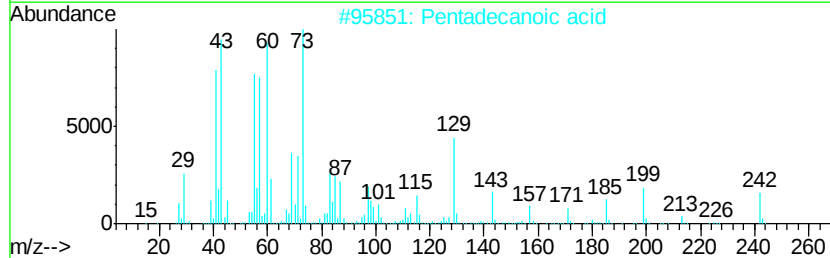
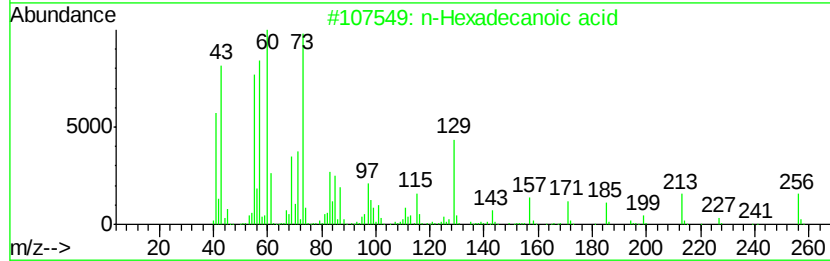
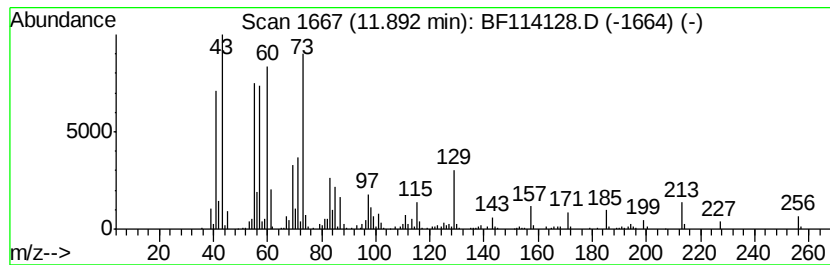
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	5.40 ng	606194	Phenanthrene-d10	11.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tridecanoic acid	214	C13H26O2	000638-53-9	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		n-Decanoic acid	172	C10H20O2	000334-48-5	53



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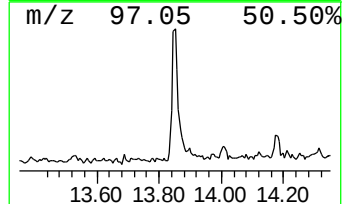
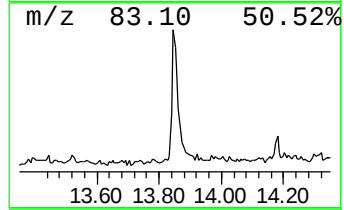
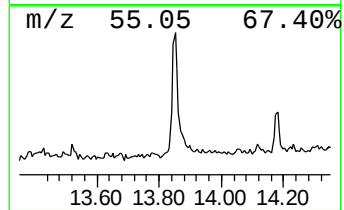
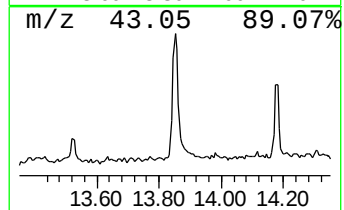
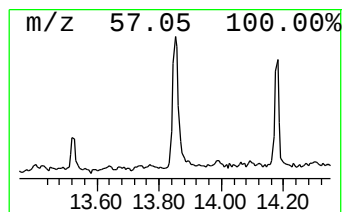
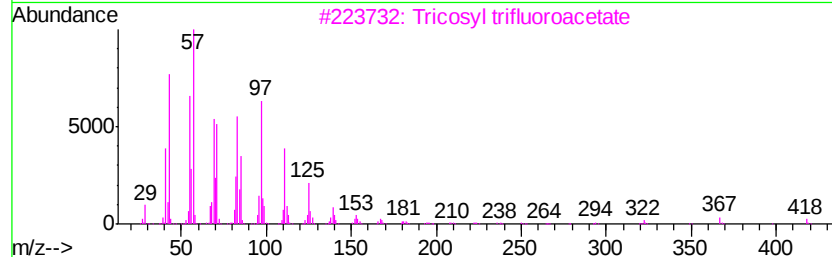
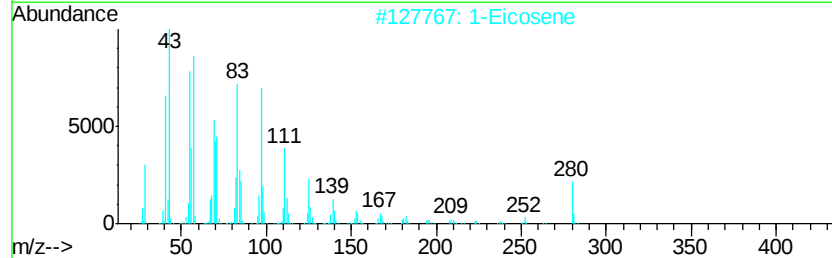
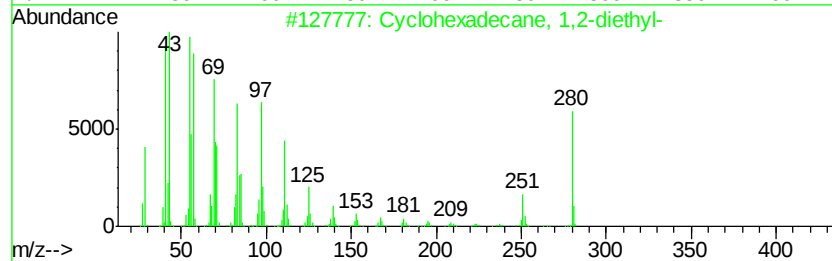
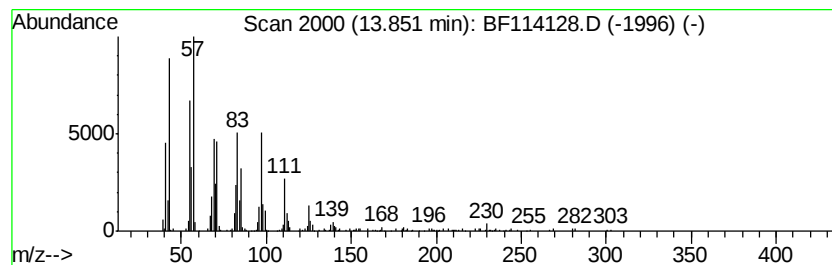
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 Peak Number 5 Cyclohexadecane, 1,2-diethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.85	5.64 ng	496096	Chrysene-d12		14.01
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	96
2	1-Eicosene	280	C20H40	003452-07-1	95
3	Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	91
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91



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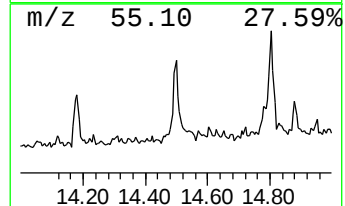
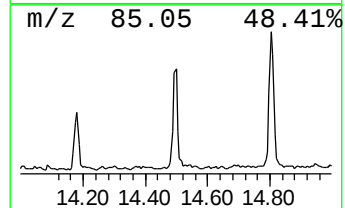
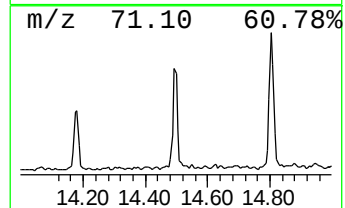
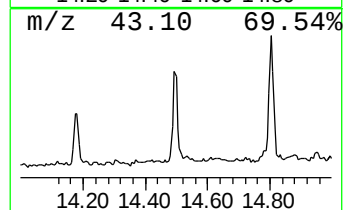
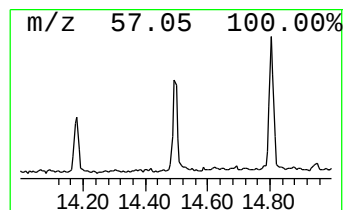
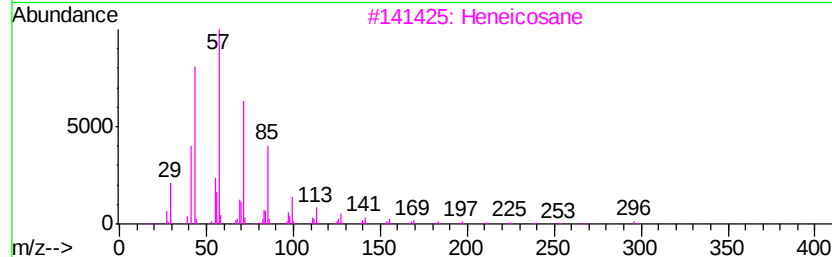
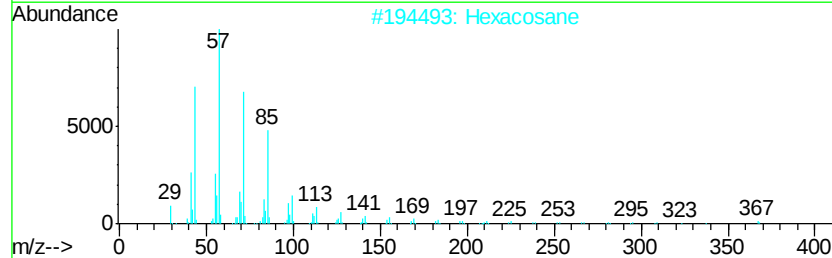
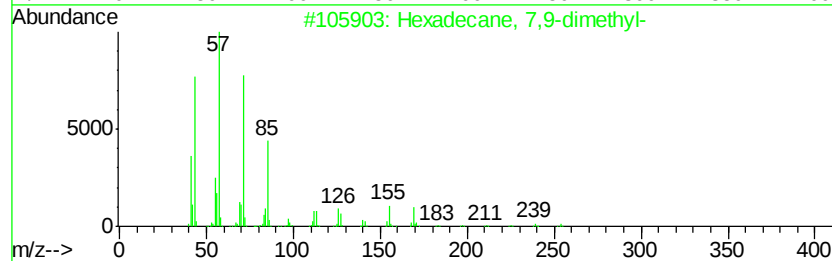
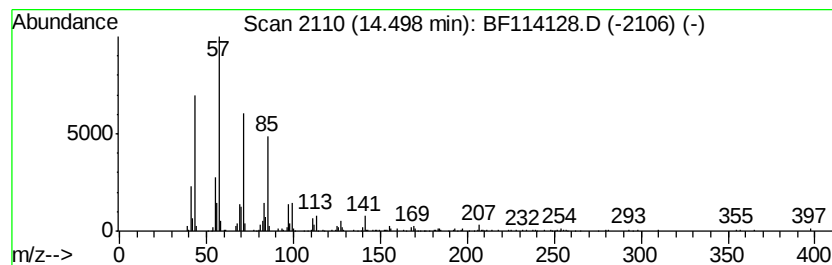
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Peak Number 6 Hexadecane, 7,9-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	2.70 ng	237076	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane, 7,9-dimethyl-	254 C18H38	021164-95-4	93
2	Hexacosane	366 C26H54	000630-01-3	90
3	Heneicosane	296 C21H44	000629-94-7	83
4	Tetracosane	338 C24H50	000646-31-1	83
5	Undecane, 2-methyl-	170 C12H26	007045-71-8	80



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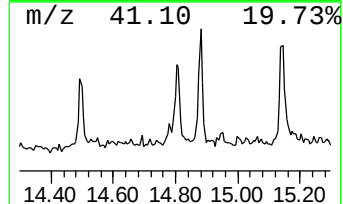
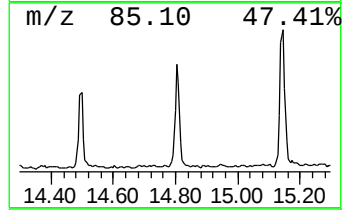
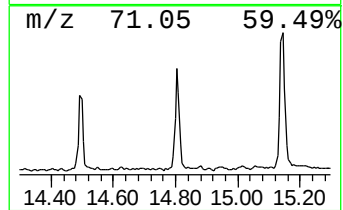
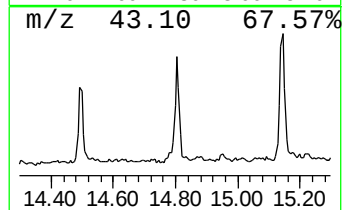
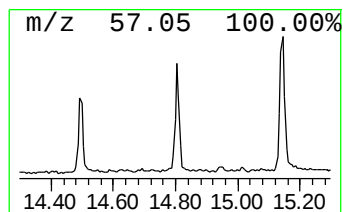
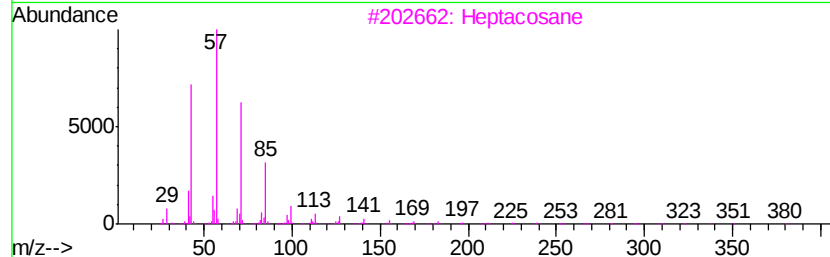
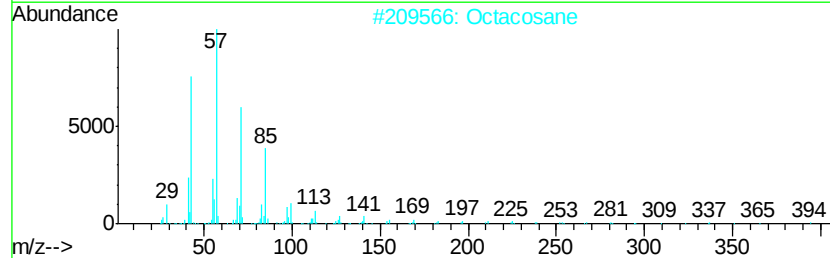
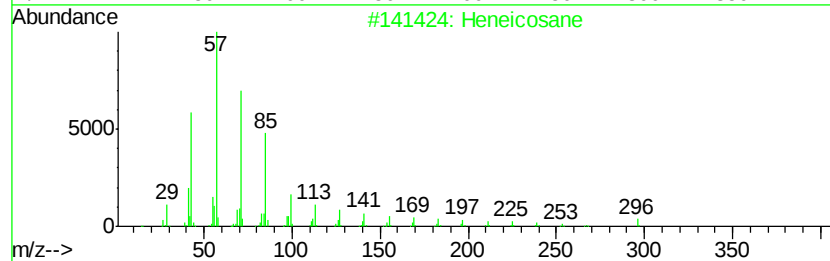
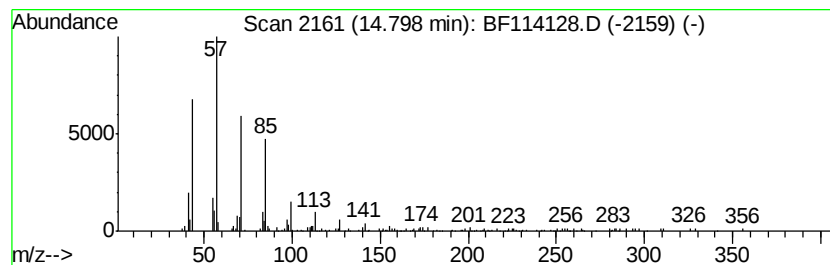
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	3.40 ng	293380	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Octadecane	254	C18H38	000593-45-3	90
5		Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Acq On : 10 May 2019 18:48
Operator : JU/SJ
Sample : K2762-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

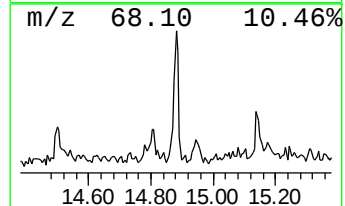
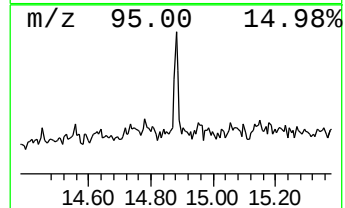
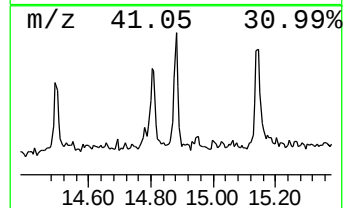
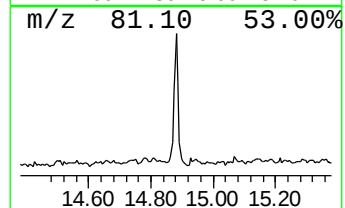
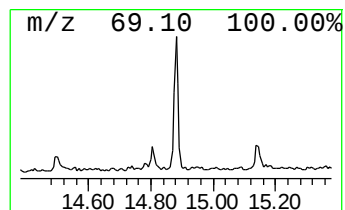
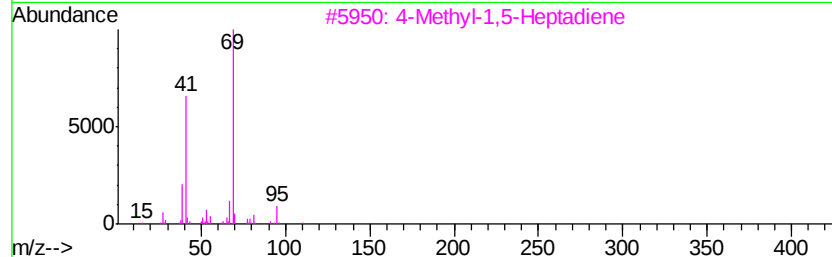
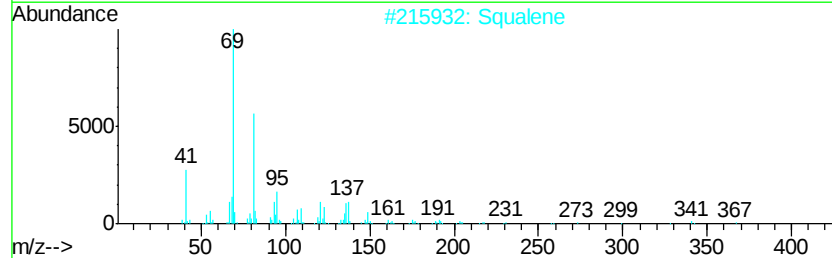
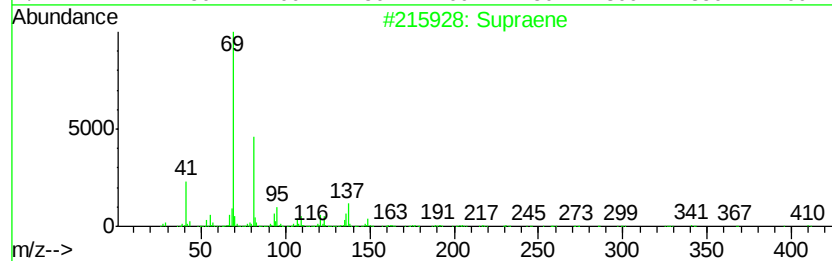
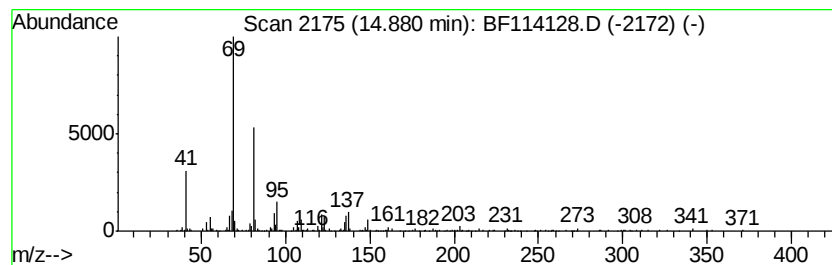
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Supraene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.88	2.68 ng	231272	Perylene-d12	15.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Supraene		410 C30H50	007683-64-9 91
2	Squalene		410 C30H50	000111-02-4 90
3	4-Methyl-1,5-Heptadiene		110 C8H14	000998-94-7 53
4	1,5-Heptadiene, 3,3,5-trimethyl-		138 C10H18	074630-29-8 53
5	1,6-Octadiene, 3,5-dimethyl-, tr...		138 C10H18	074630-87-8 53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Instrument :
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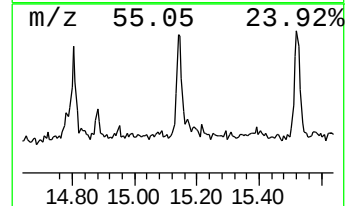
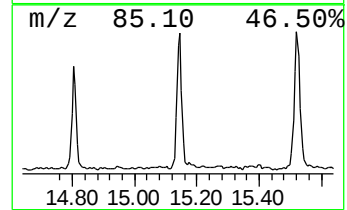
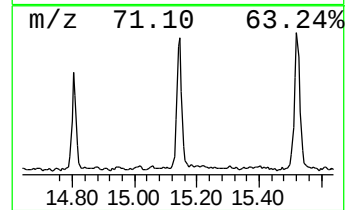
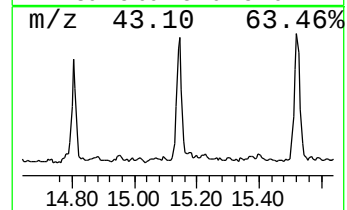
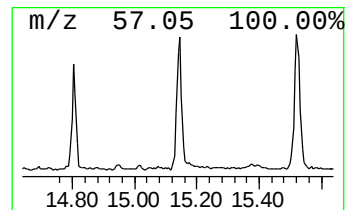
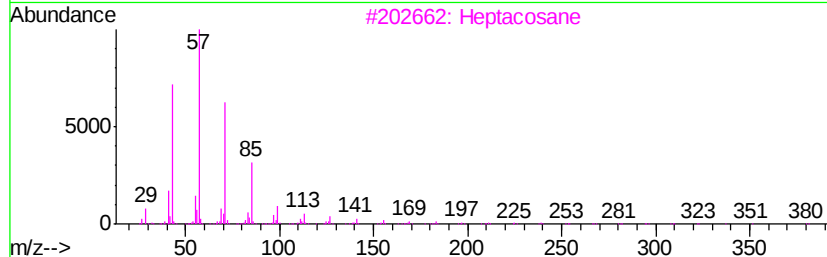
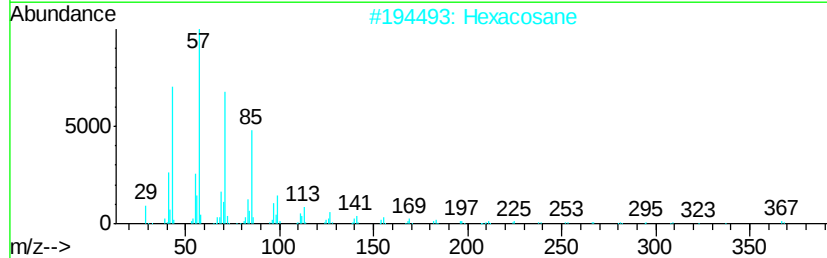
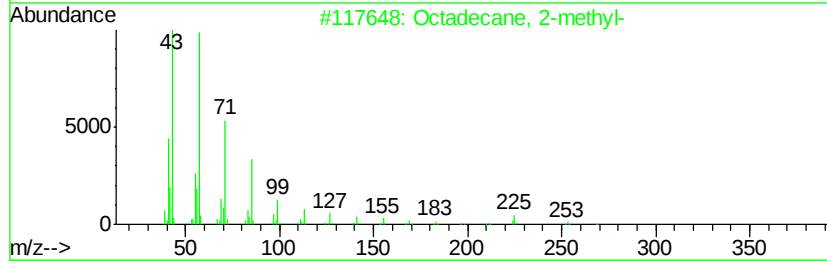
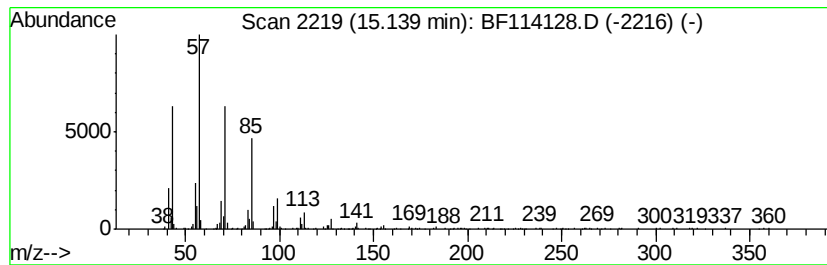
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 9 Octadecane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	5.38 ng	464610	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heptacosane	380	C27H56	000593-49-7	87
4		Octadecane	254	C18H38	000593-45-3	87
5		Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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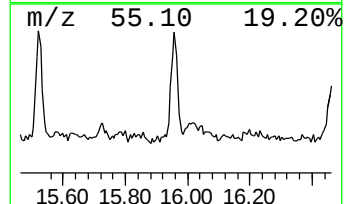
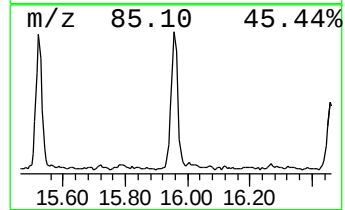
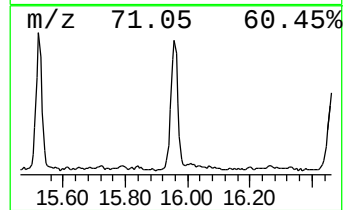
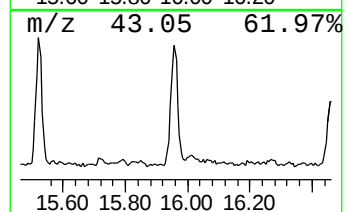
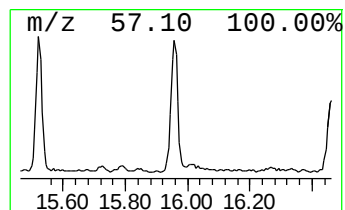
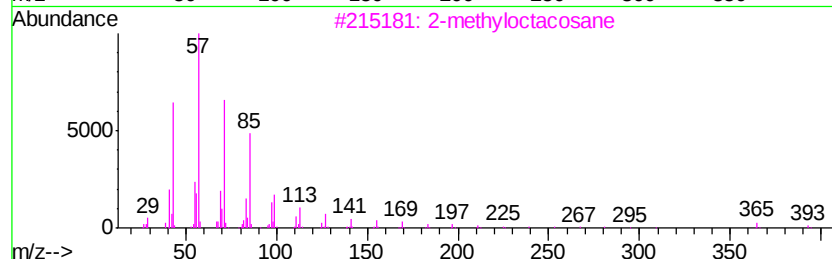
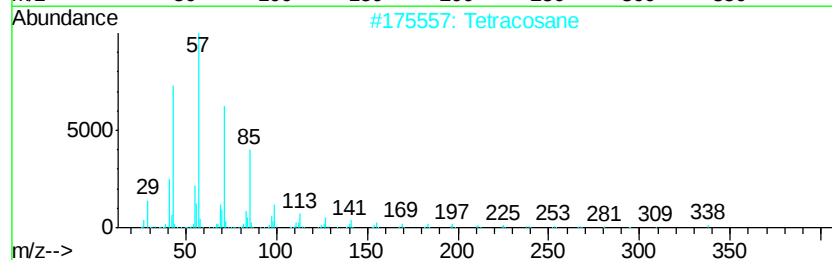
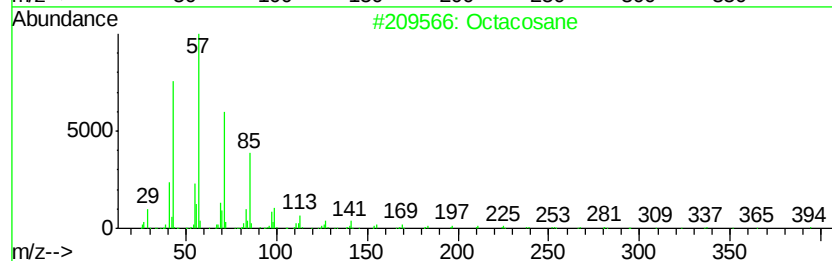
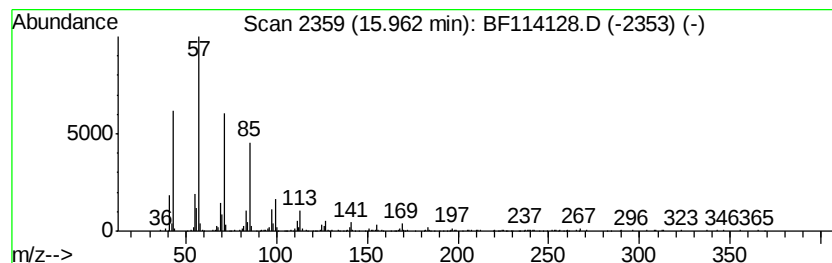
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Octacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.96	7.51 ng	648645	Perylene-d12	15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			2-methyloctacosane	408	C29H60	1000376-72-8	90
4			Octadecane	254	C18H38	000593-45-3	87
5			Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
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Instrument :
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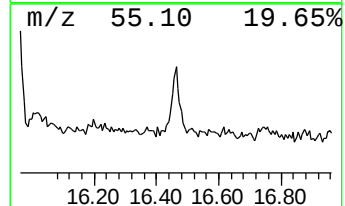
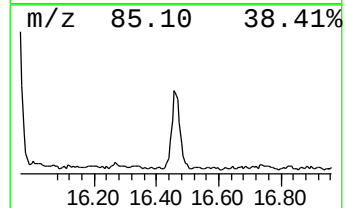
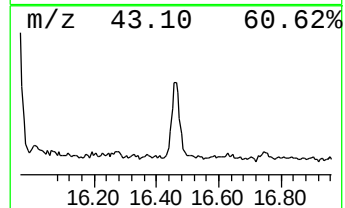
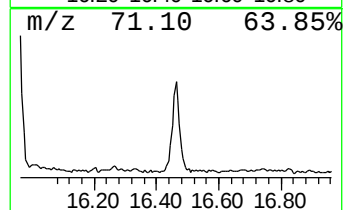
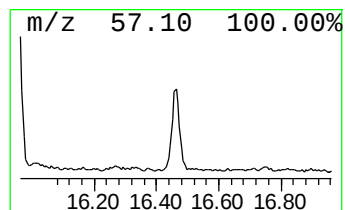
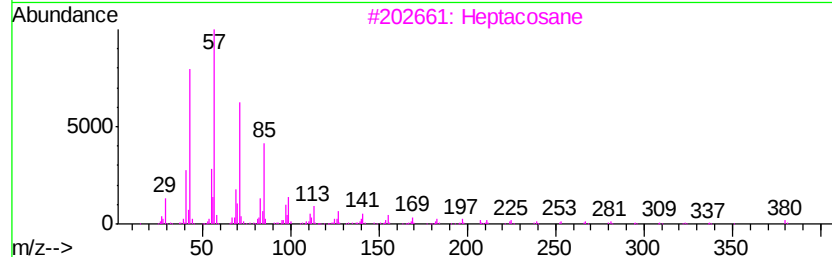
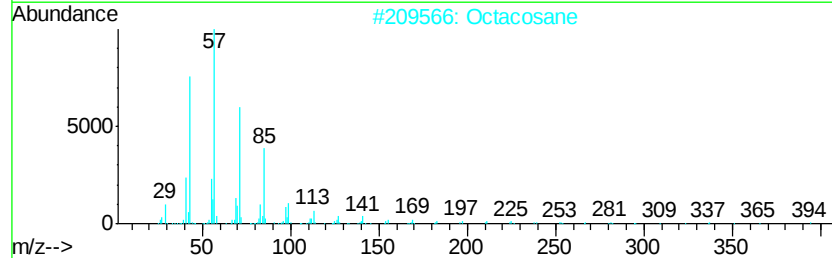
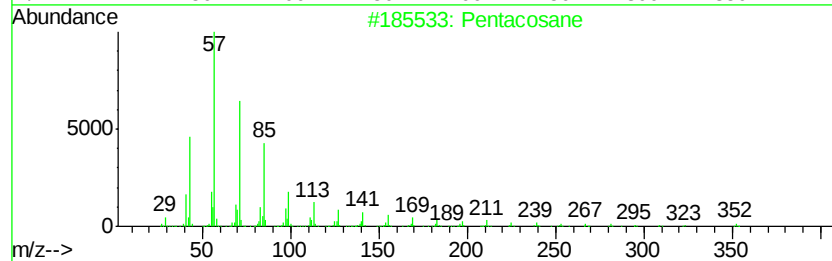
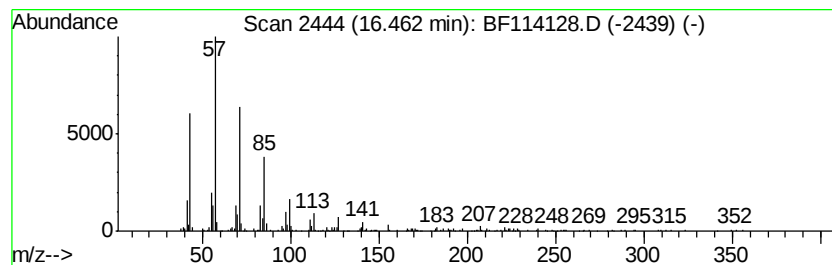
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Pentacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	4.50 ng	388994	Perylene-d12	15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	95
2			Octacosane	394	C28H58	000630-02-4	90
3			Heptacosane	380	C27H56	000593-49-7	90
4			Hexacosane	366	C26H54	000630-01-3	90
5			Hentriacontane	437	C31H64	000630-04-6	90



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Data File : BF114128.D
Acq On : 10 May 2019 18:48
Operator : JU/SJ
Sample : K2762-10
Misc :
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Instrument :
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ClientSampleId :
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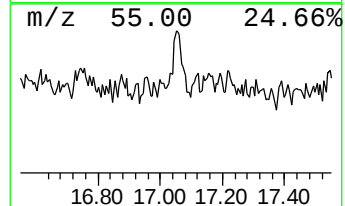
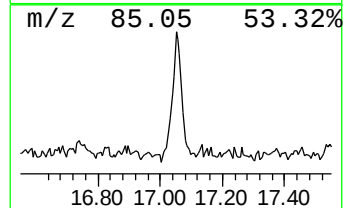
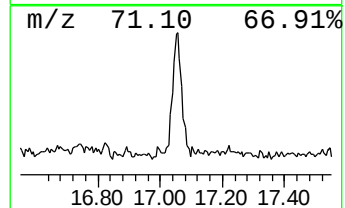
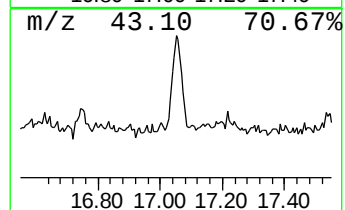
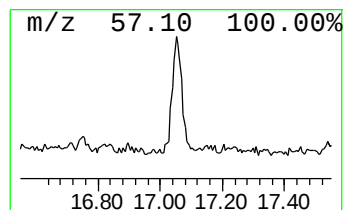
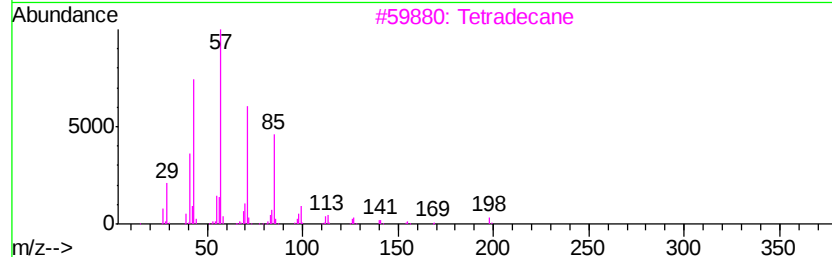
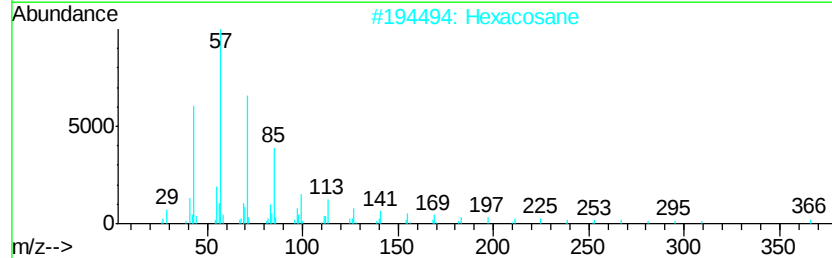
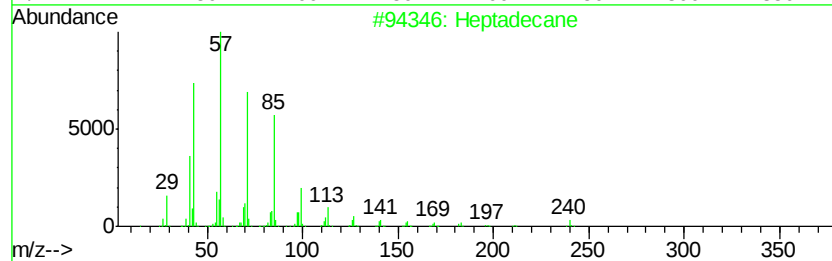
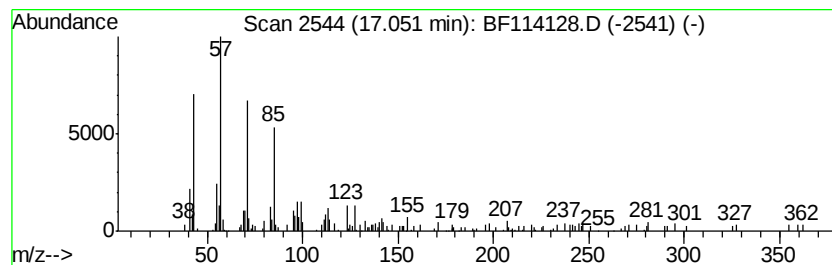
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.05	2.25 ng	194147	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Hexacosane	366	C26H54	000630-01-3	70
3		Tetradecane	198	C14H30	000629-59-4	64
4		Heptacosane	380	C27H56	000593-49-7	62
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	62



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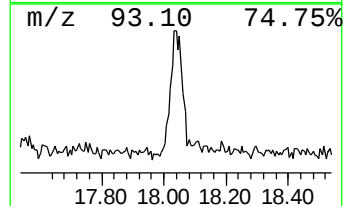
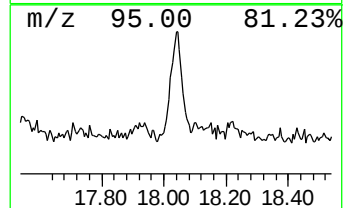
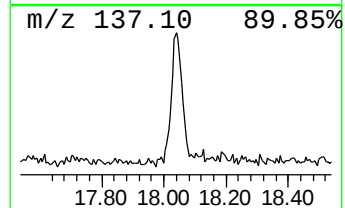
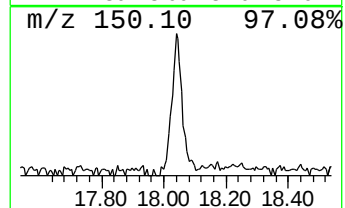
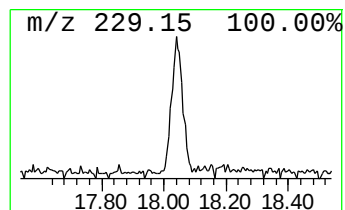
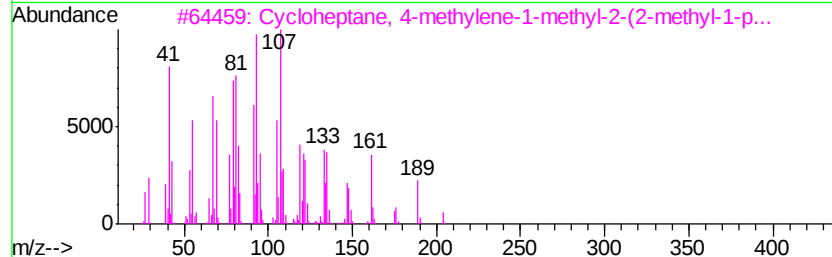
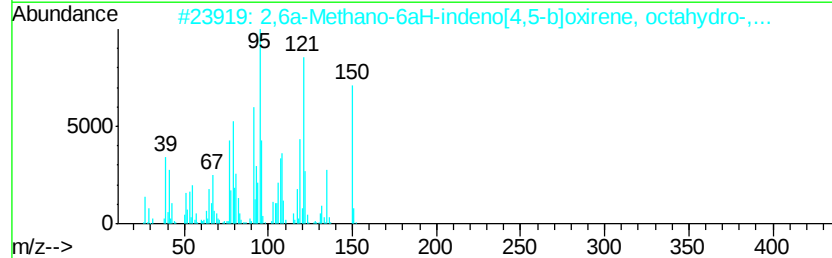
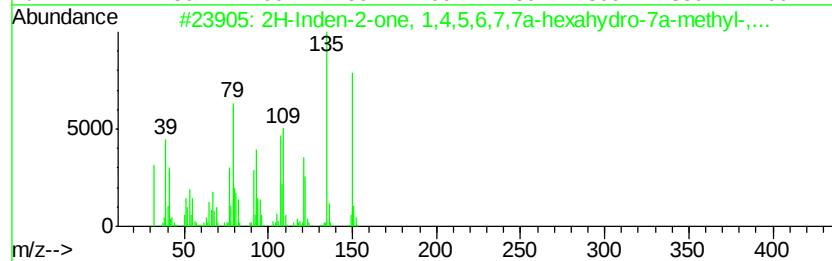
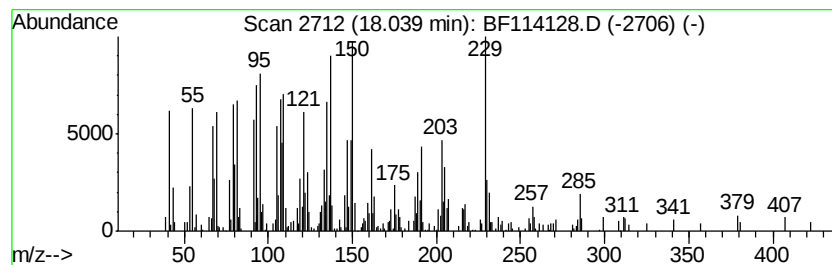
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown18.04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	6.11 ng	527449	Perylene-d12	15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Inden-2-one, 1,4,5,6,7,7a-hex...	150	C10H14O	054725-16-5	47
2			2,6a-Methano-6aH-indeno[4,5-b]ox...	150	C10H14O	016489-32-0	47
3			Cycloheptane, 4-methylene-1-meth...	204	C15H24	1000159-38-5	44
4			Cyclohexane, 1,1,2-trimethyl-3,5...	206	C15H26	062337-97-7	38
5			2-Cyclopenten-1-one, 2-(2-buteny...	150	C10H14O	017190-71-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051019\
 Data File : BF114128.D
 Acq On : 10 May 2019 18:48
 Operator : JU/SJ
 Sample : K2762-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P021-SS002-0206-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.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...	2.15	18.0	ng	1713450	1	6.85	1908040	20.0
unknown6.62	6.62	75.0	ng	7151950	1	6.85	1908040	20.0
n-Hexadecanoic acid	11.89	5.4	ng	606194	4	11.37	2245620	20.0
Cyclohexadecane, ...	13.85	5.6	ng	496096	5	14.01	1759060	20.0
Hexadecane, 7,9-d...	14.50	2.7	ng	237076	5	14.01	1759060	20.0
Heneicosane	14.80	3.4	ng	293380	6	15.49	1727030	20.0
Supraene	14.88	2.7	ng	231272	6	15.49	1727030	20.0
Octadecane, 2-met...	15.14	5.4	ng	464610	6	15.49	1727030	20.0
Octacosane	15.96	7.5	ng	648645	6	15.49	1727030	20.0
Pentacosane	16.46	4.5	ng	388994	6	15.49	1727030	20.0
Heptadecane	17.05	2.3	ng	194147	6	15.49	1727030	20.0
unknown18.04	18.04	6.1	ng	527449	6	15.49	1727030	20.0