

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111286.D
 Acq On : 11 Dec 2018 19:36
 Operator : JU/SJ
 Sample : J6157-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMPOSITE-2-DISSOLVED

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-BF120818.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	2.828	120	126	132	rVB	242492	390186	7.53%	0.793%
2	3.793	286	290	299	rBV2	38059	66426	1.28%	0.135%
3	4.034	327	331	336	rVB3	60322	73851	1.43%	0.150%
4	4.987	484	493	498	rBV2	255998	317497	6.13%	0.645%
5	5.404	560	564	568	rBV	2735059	2555255	49.32%	5.191%
6	5.922	649	652	655	rVB	65847	57464	1.11%	0.117%
7	6.422	733	737	749	rBV2	1665776	2112272	40.77%	4.291%
8	6.563	757	761	764	rBV	4811952	4086011	78.87%	8.300%
9	6.792	797	800	804	rBV	1577046	1448289	27.96%	2.942%
10	6.863	809	812	817	rBV	62629	62015	1.20%	0.126%
11	6.951	823	827	831	rBV	4424963	3820219	73.74%	7.760%
12	7.357	890	896	899	rBV	3128507	3100638	59.85%	6.299%
13	8.081	1014	1019	1022	rBV	1909668	1802119	34.78%	3.661%
14	9.157	1197	1202	1205	rBV	5890973	5180766	100.00%	10.524%
15	9.545	1264	1268	1272	rBV	1027694	824382	15.91%	1.675%
16	9.833	1313	1317	1321	rVB	2817537	2392002	46.17%	4.859%
17	10.251	1384	1388	1390	rBV	282355	229640	4.43%	0.466%
18	10.622	1446	1451	1454	rBV	3205974	3129461	60.41%	6.357%
19	11.316	1565	1569	1572	rBV	2827590	2259629	43.62%	4.590%
20	11.710	1632	1636	1639	rBV2	157167	137020	2.64%	0.278%
21	11.857	1654	1661	1670	rBV	2330521	2341636	45.20%	4.757%
22	12.539	1774	1777	1778	rBV2	76747	65665	1.27%	0.133%
23	12.557	1778	1780	1787	rVB2	64587	73177	1.41%	0.149%
24	12.639	1790	1794	1804	rVV	1422737	1379572	26.63%	2.802%
25	12.733	1807	1810	1814	rVB	111515	106039	2.05%	0.215%
26	12.904	1834	1839	1843	rBV	4916520	5089220	98.23%	10.338%
27	13.810	1989	1993	2000	rBV2	870152	1119061	21.60%	2.273%
28	13.957	2013	2018	2021	rVB	2228088	1991130	38.43%	4.045%
29	14.133	2046	2048	2055	rBV	133534	178557	3.45%	0.363%
30	14.298	2073	2076	2082	rBV8	33940	68536	1.32%	0.139%
31	14.445	2098	2101	2112	rVB2	196597	314498	6.07%	0.639%
32	14.745	2149	2152	2157	rVB	132016	135809	2.62%	0.276%
33	15.080	2205	2209	2212	rBV	183273	189804	3.66%	0.386%
34	15.398	2258	2263	2268	rBV	757063	941594	18.17%	1.913%

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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	15.451	2268	2272	2276	rVB	176049	199801	3.86%	0.406%
36	15.874	2340	2344	2349	rBV	151861	225203	4.35%	0.457%
37	15.927	2350	2353	2360	rVV3	57582	99548	1.92%	0.202%
38	16.368	2423	2428	2436	rBV3	123557	226288	4.37%	0.460%
39	16.609	2464	2469	2478	rVB6	71136	154130	2.98%	0.313%
40	16.951	2521	2527	2532	rBV2	77521	148750	2.87%	0.302%
41	17.627	2637	2642	2650	rVB5	57618	134342	2.59%	0.273%

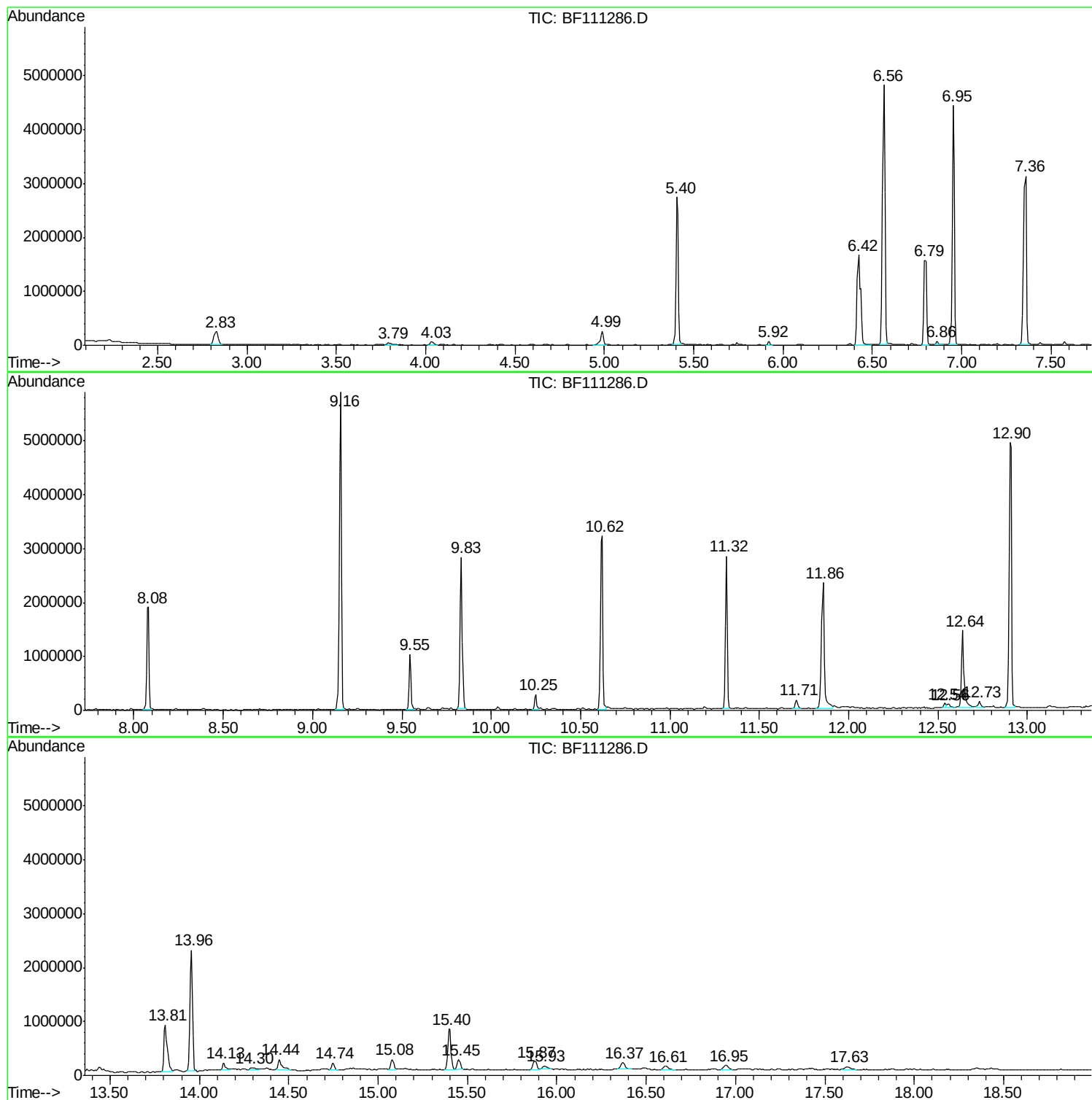
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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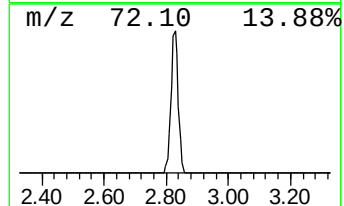
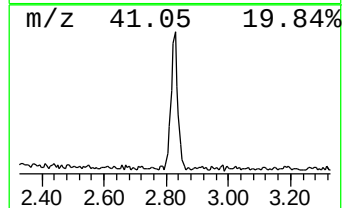
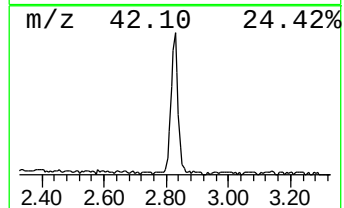
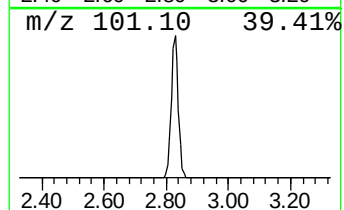
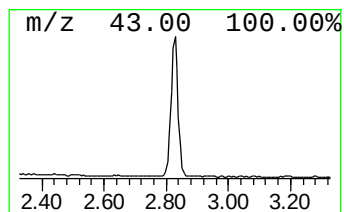
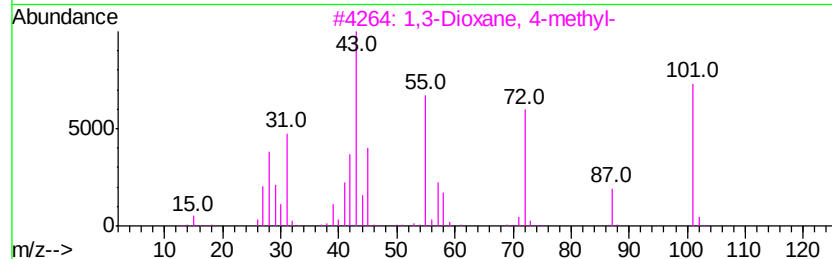
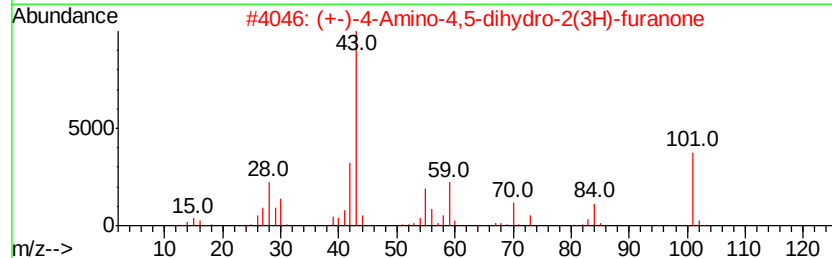
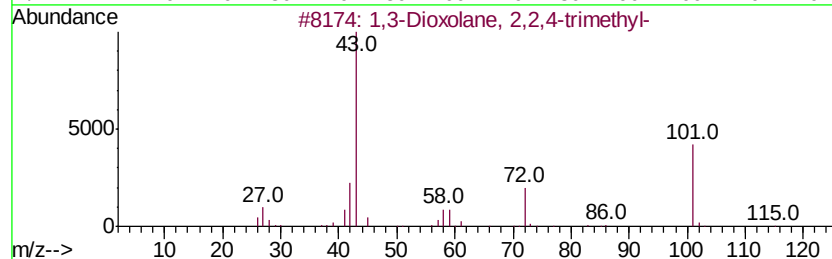
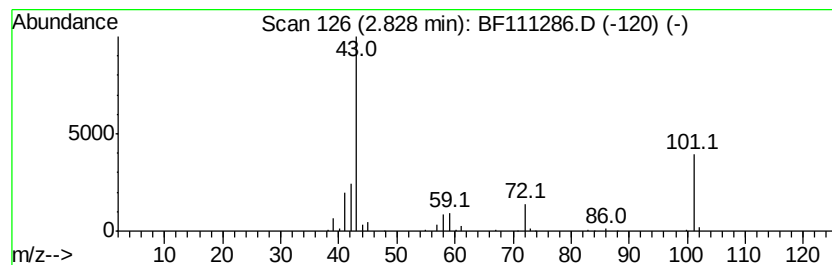
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	5.39 ng	390186	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	78
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	38
3		1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	36
4		Furazan-3-ol, 4-amino-	101	C2H3N3O2	159013-90-8	9
5		Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	9



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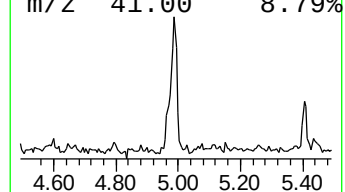
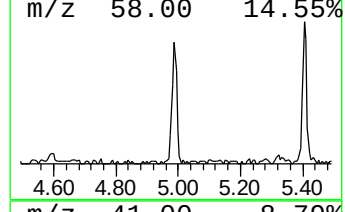
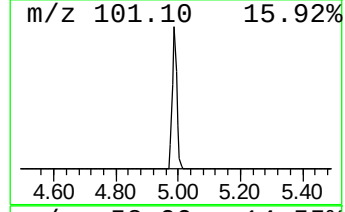
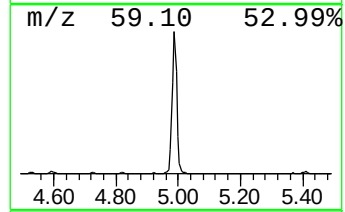
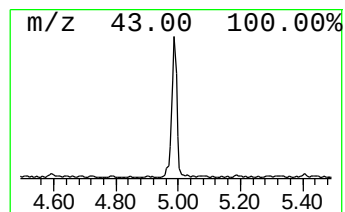
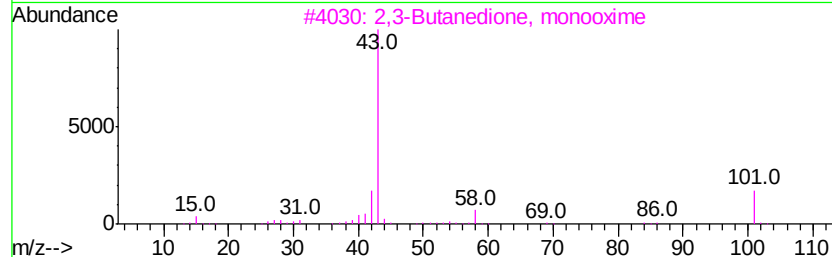
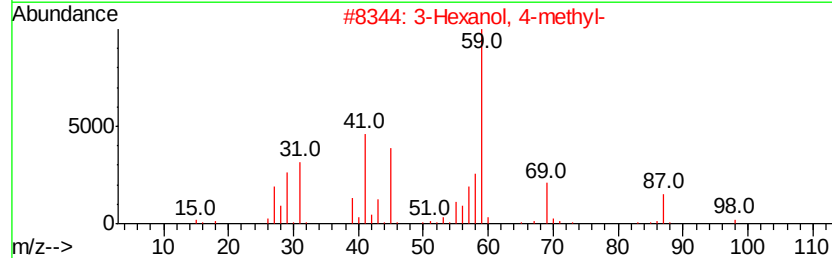
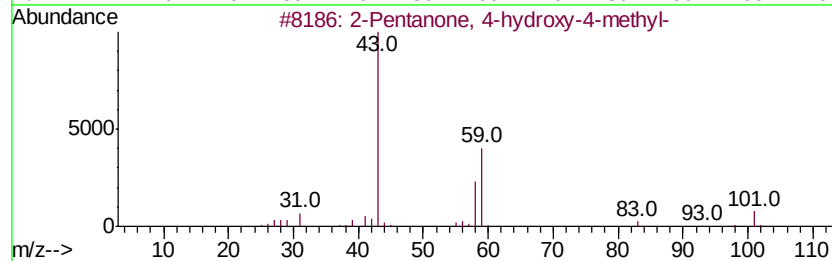
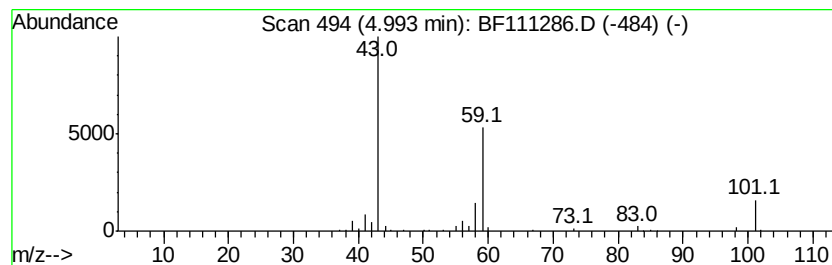
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	4.38 ng	317497	1,4-Dichlorobenzene-d4	6.80
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116 C6H12O2	000123-42-2	45
2	3-Hexanol, 4-methyl-	116 C7H16O	000615-29-2	33
3	2,3-Butanedione, monooxime	101 C4H7NO2	000057-71-6	27
4	Acetic acid, cyano-, 1,1-dimethyl...	141 C7H11NO2	001116-98-9	25
5	Formamide, N,N-diethyl-	101 C5H11NO	000617-84-5	9



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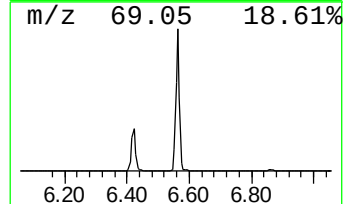
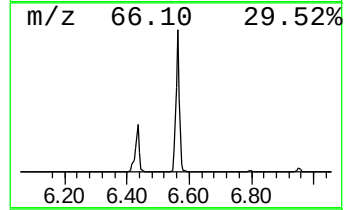
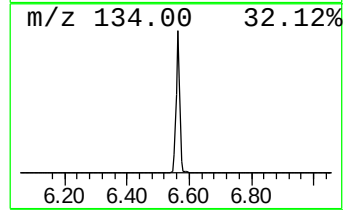
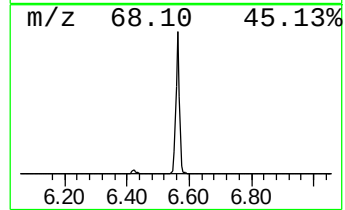
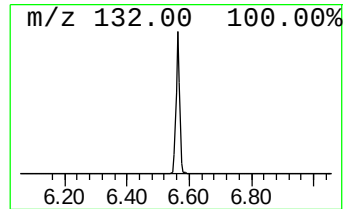
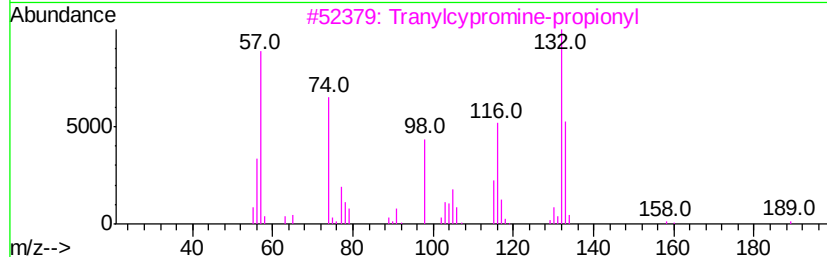
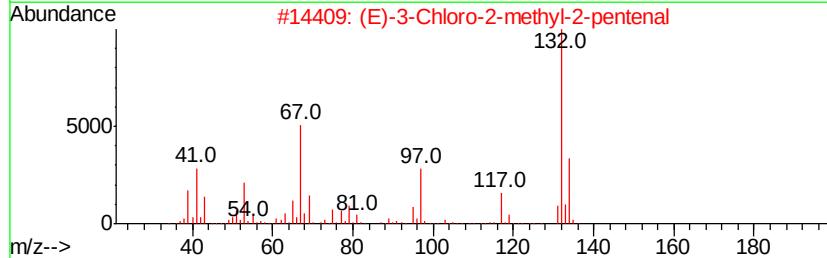
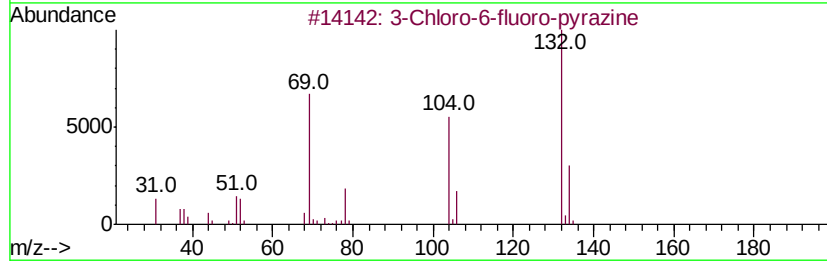
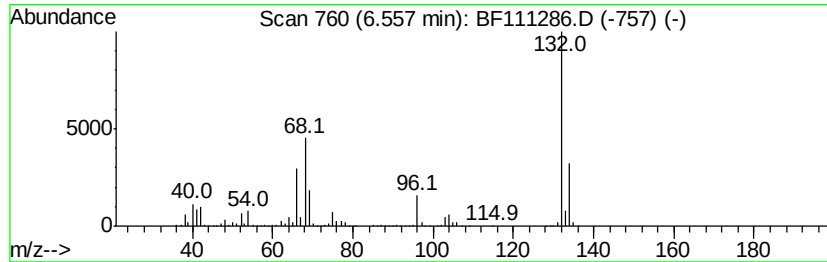
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	56.43 ng	4086010	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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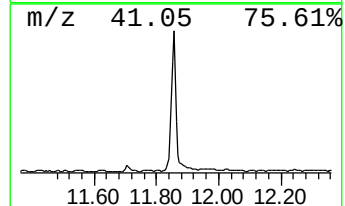
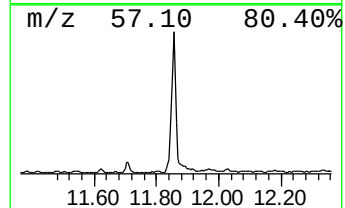
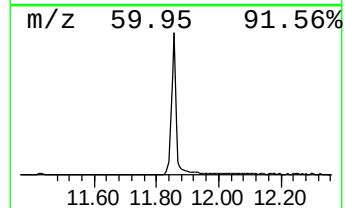
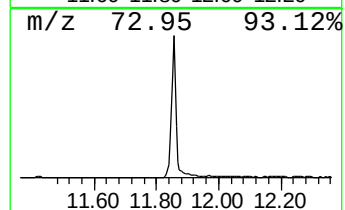
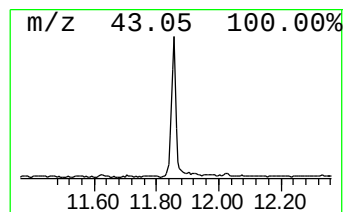
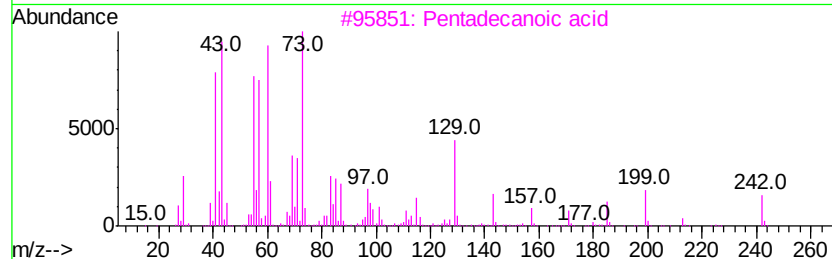
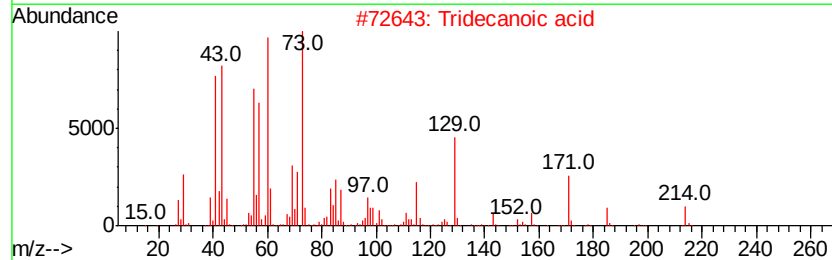
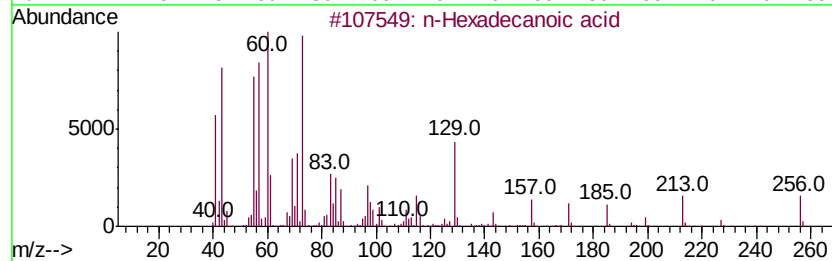
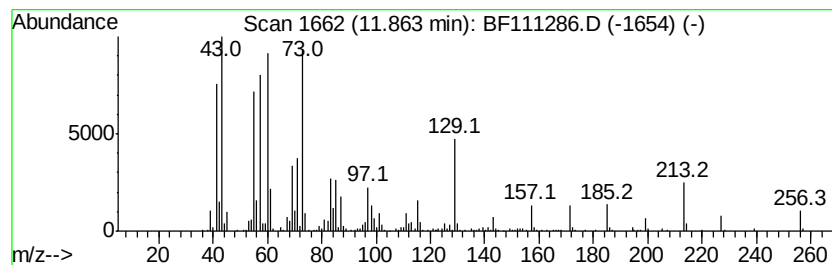
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	20.73 ng	2341640	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Undecanoic acid	186	C11H22O2	000112-37-8	72



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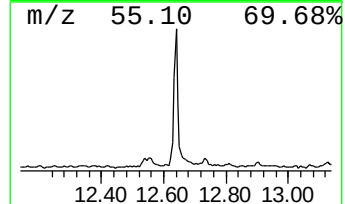
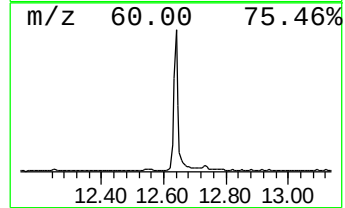
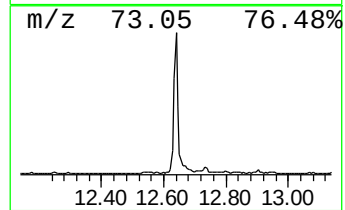
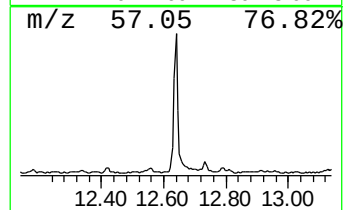
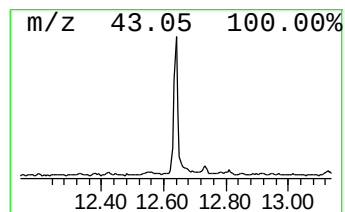
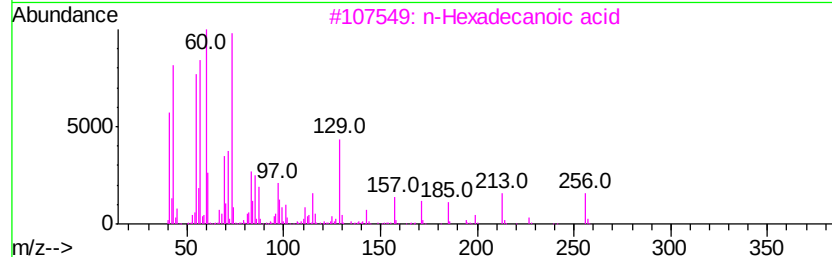
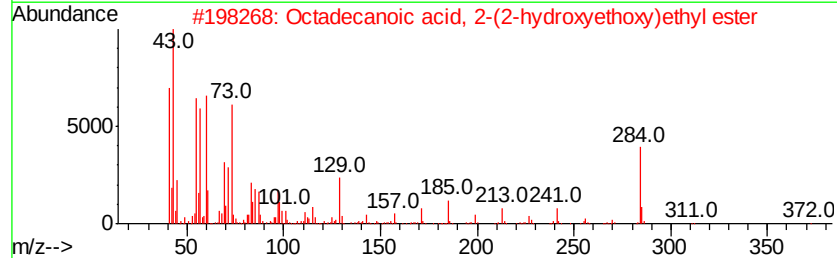
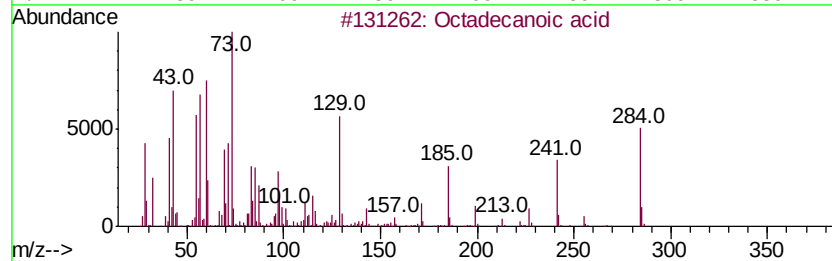
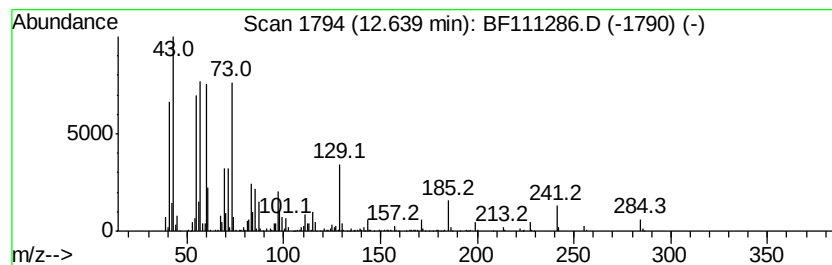
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	13.86 ng	1379570	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	91
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5		Tridecanoic acid	214	C13H26O2	000638-53-9	74



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Acq On : 11 Dec 2018 19:36
Operator : JU/SJ
Sample : J6157-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

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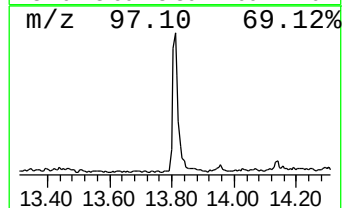
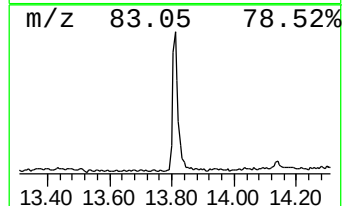
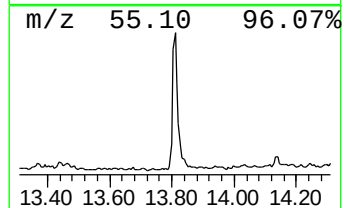
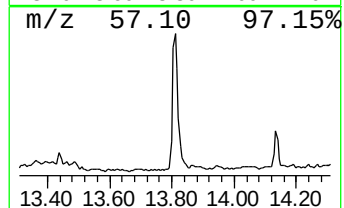
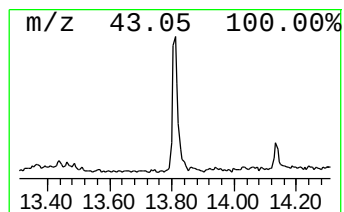
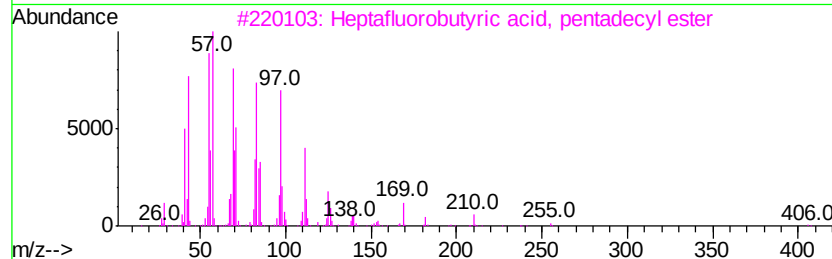
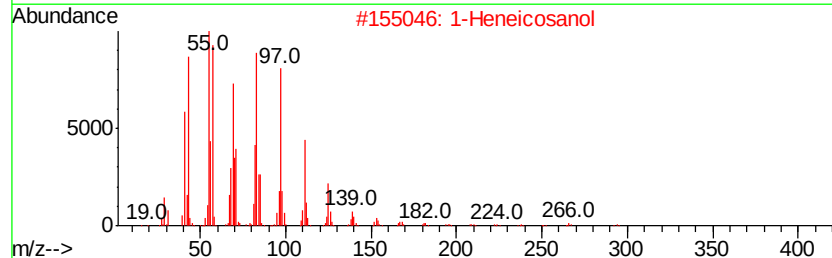
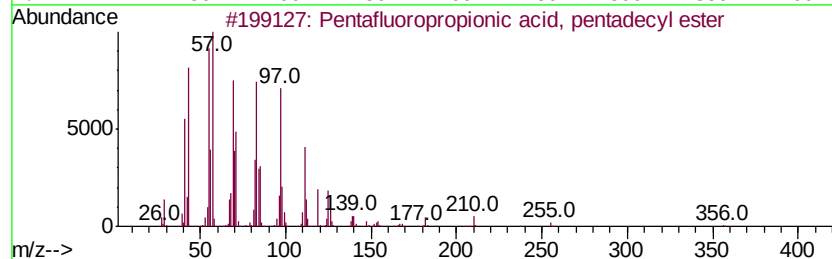
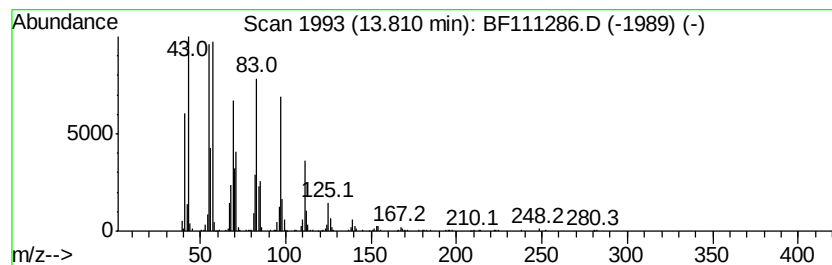
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentafluoropropionic acid, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	11.24 ng	1119060	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
2		1-Heneicosanol	312	C21H44O	015594-90-8	94
3		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4		1-Docosene	308	C22H44	001599-67-3	94
5		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94



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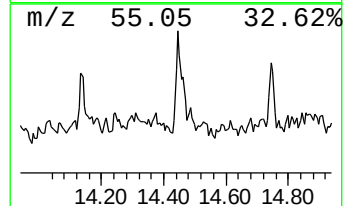
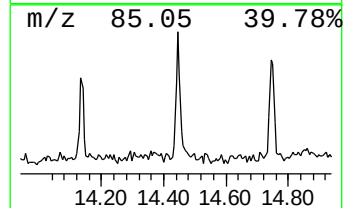
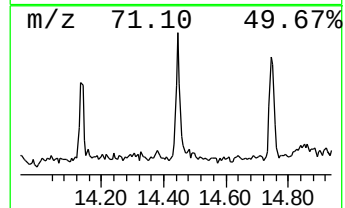
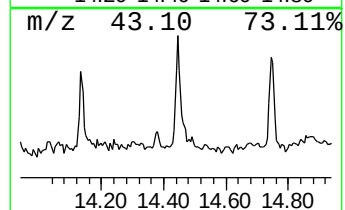
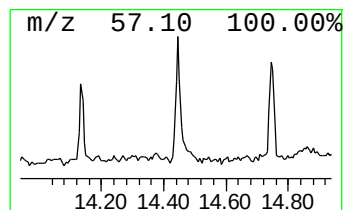
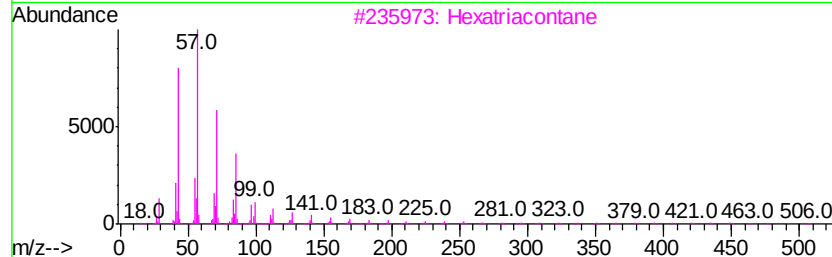
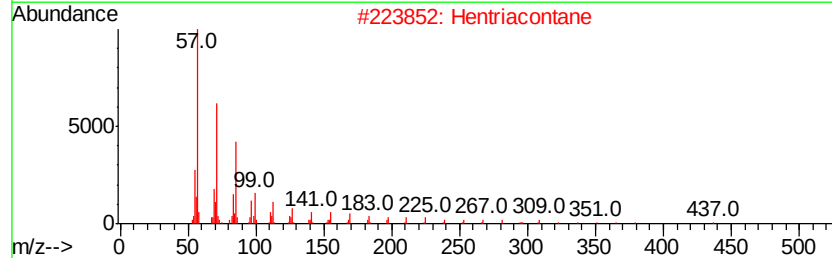
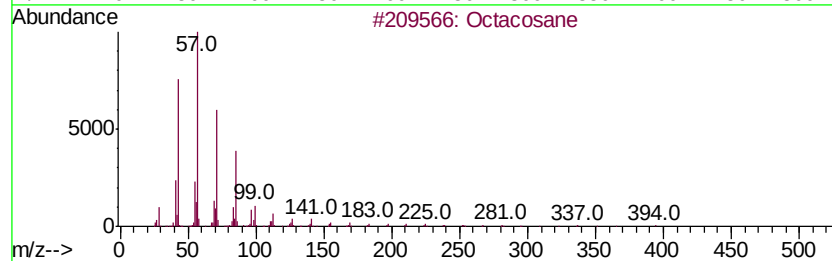
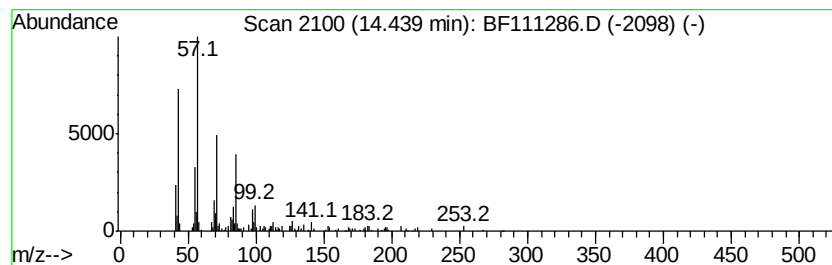
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Octacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
14.44	3.16 ng	314498	Chrysene-d12		13.96		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Hentriacontane	437	C31H64	000630-04-6	87
3			Hexatriacontane	507	C36H74	000630-06-8	86
4			Heneicosane	296	C21H44	000629-94-7	83
5			Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
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 Acq On : 11 Dec 2018 19:36
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Instrument :
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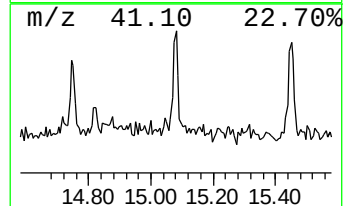
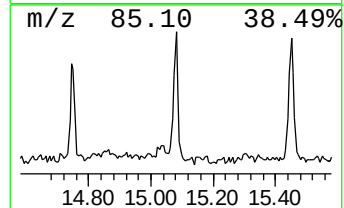
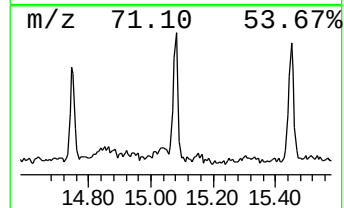
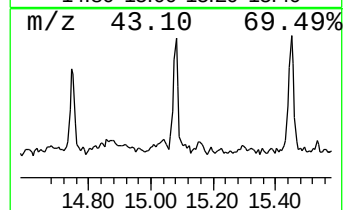
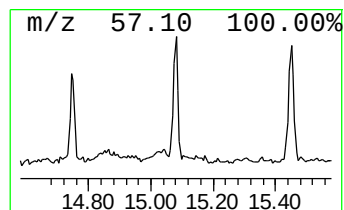
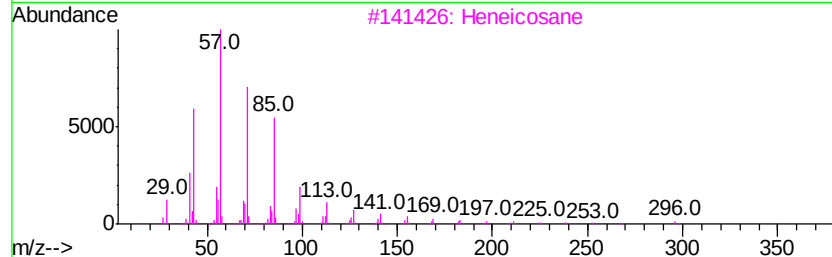
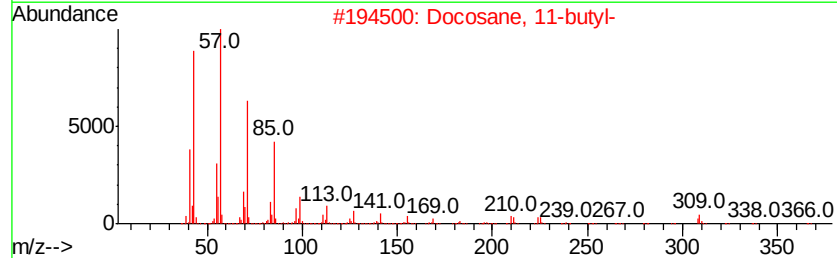
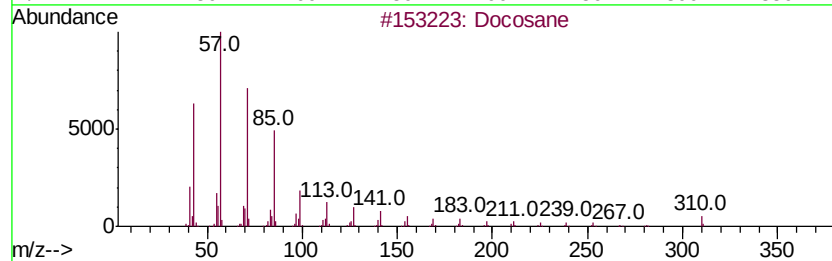
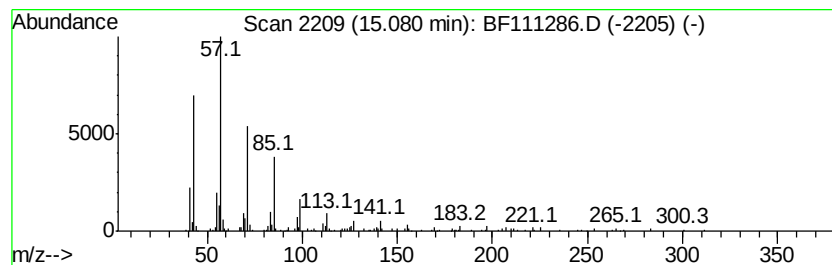
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Docosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.08	4.03 ng	189804	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	90
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Octadecane	254	C18H38	000593-45-3	90
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
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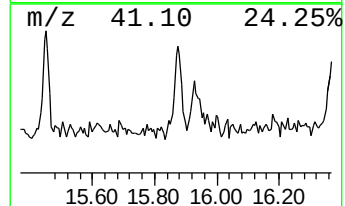
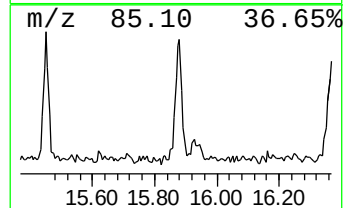
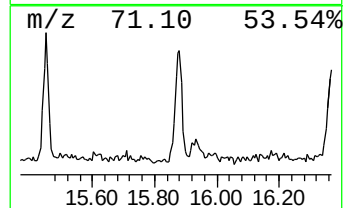
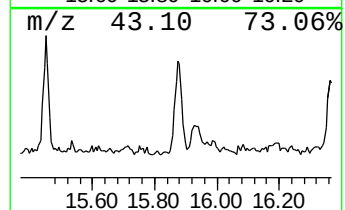
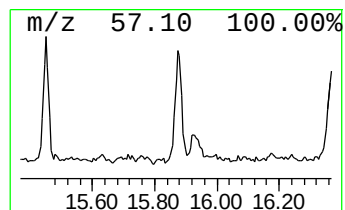
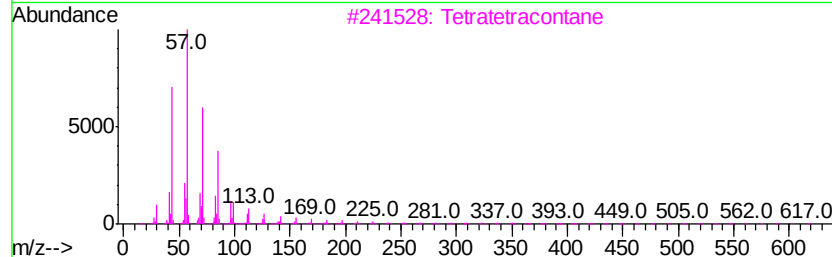
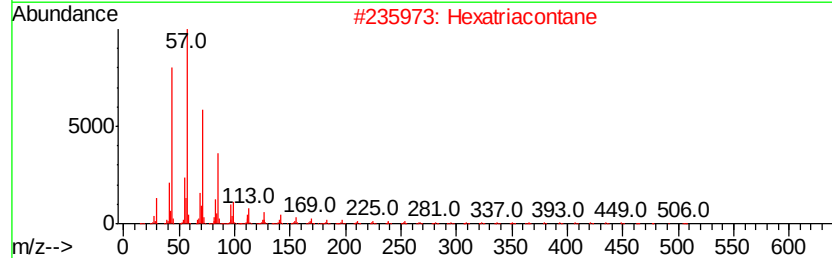
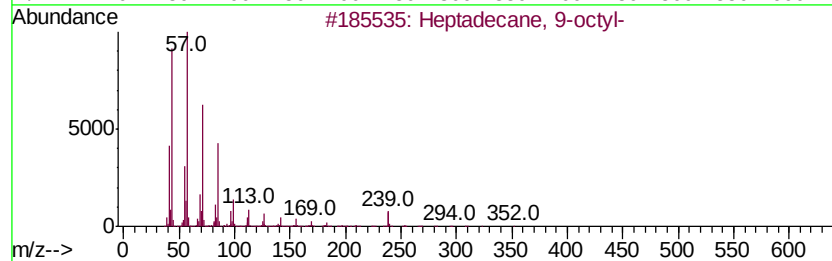
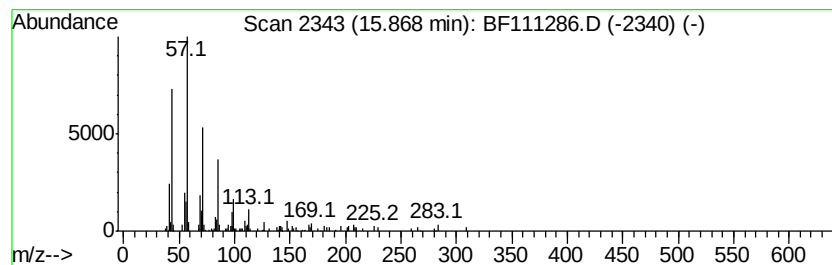
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Heptadecane, 9-octyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	4.78 ng	225203	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
2		Hexatriacontane	507	C36H74	000630-06-8	90
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
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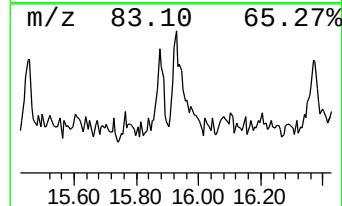
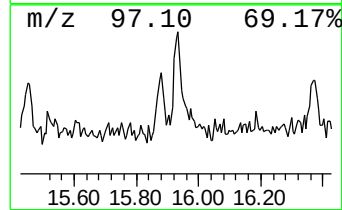
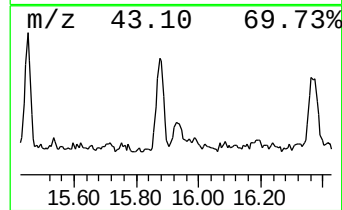
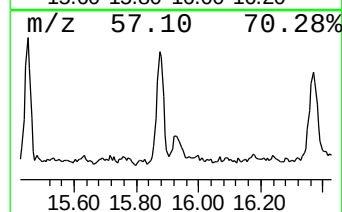
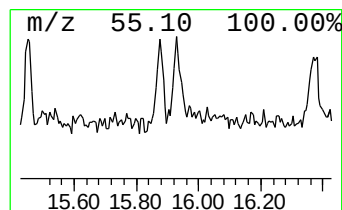
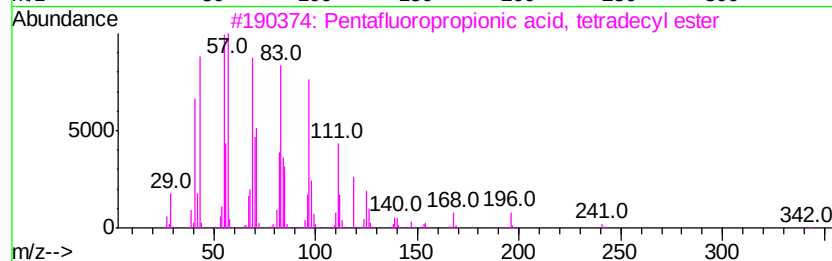
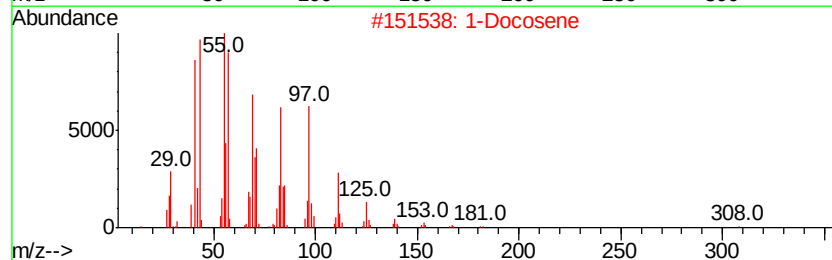
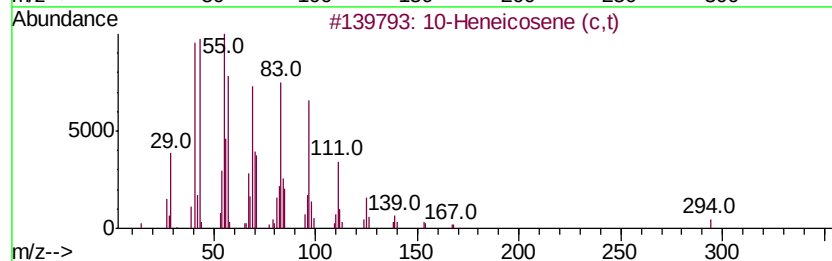
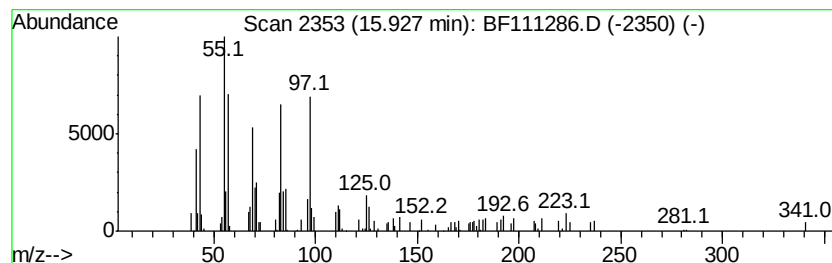
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 10-Heneicosene (c,t) Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	2.11 ng	99548	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Heneicosene (c,t)	294	C21H42	095008-11-0	64
2		1-Docosene	308	C22H44	001599-67-3	64
3		Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	58
4		11-Methyldodecanol	200	C13H28O	085763-57-1	52
5		1-Tridecene	182	C13H26	002437-56-1	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
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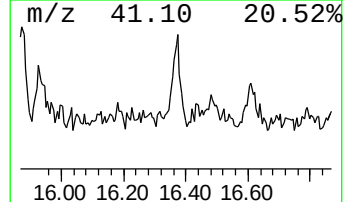
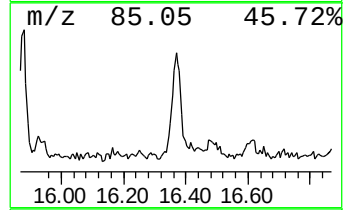
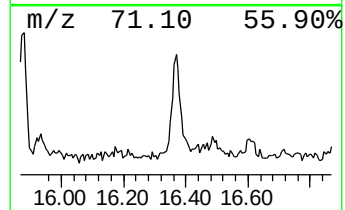
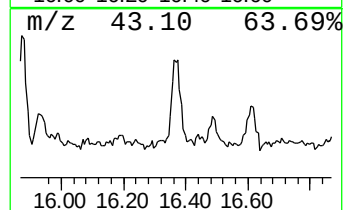
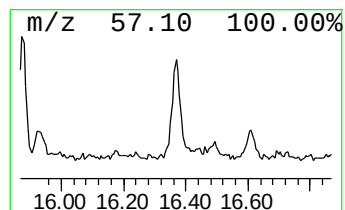
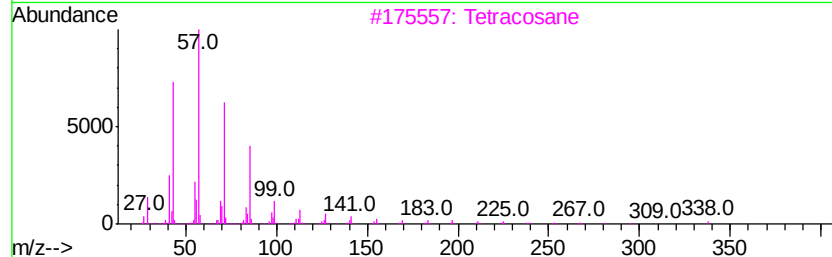
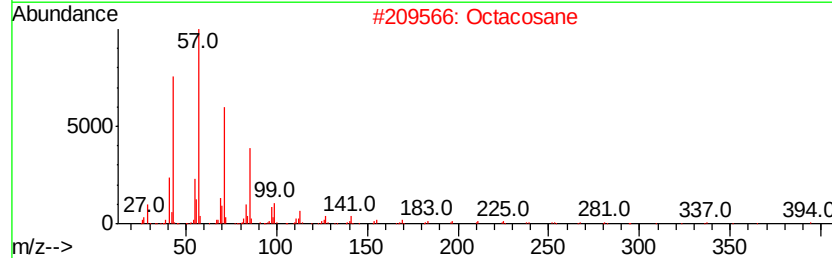
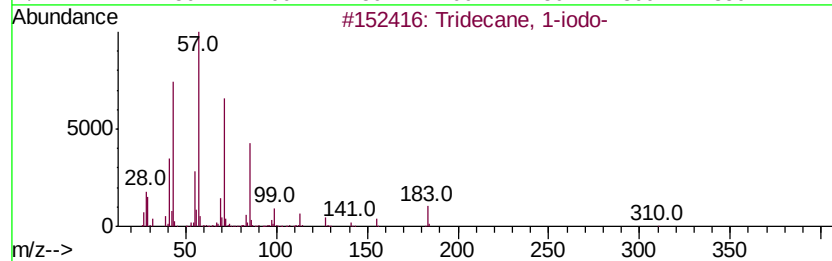
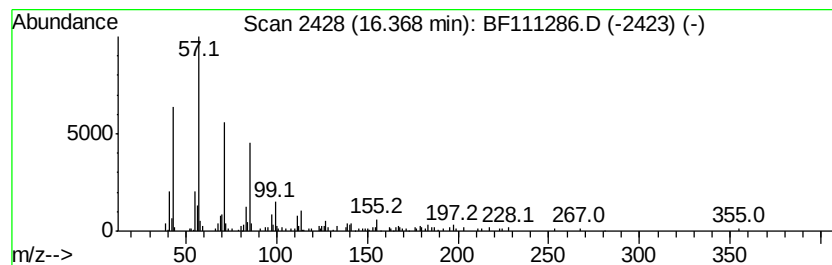
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tridecane, 1-iodo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	4.81 ng	226288	Perylene-d12	15.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Tridecane, 1-iodo-	310 C13H27I	035599-77-0	90
2	Octacosane	394 C28H58	000630-02-4	87
3	Tetracosane	338 C24H50	000646-31-1	87
4	Nonadecane	268 C19H40	000629-92-5	87
5	Docosane	310 C22H46	000629-97-0	86



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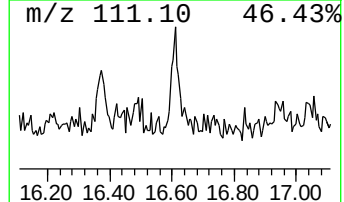
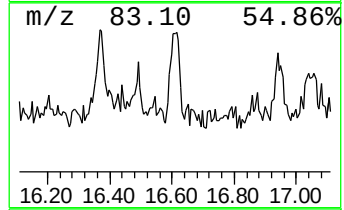
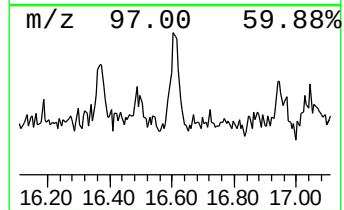
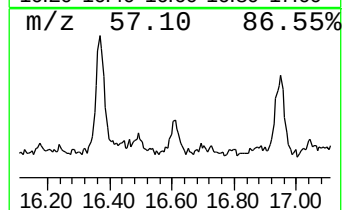
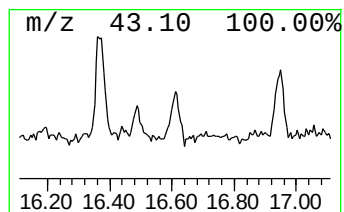
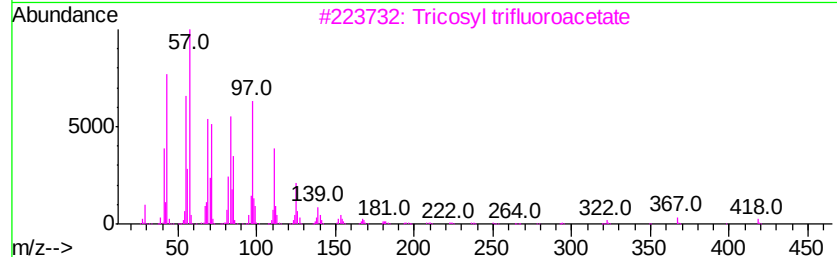
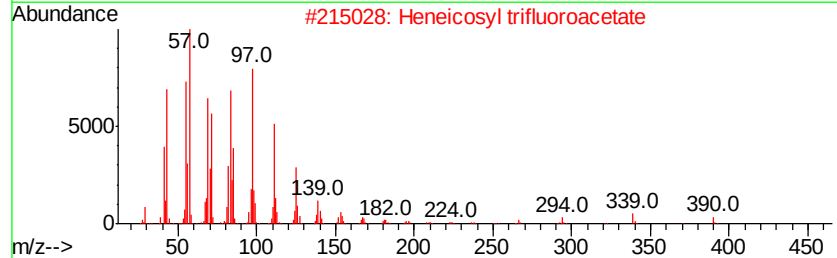
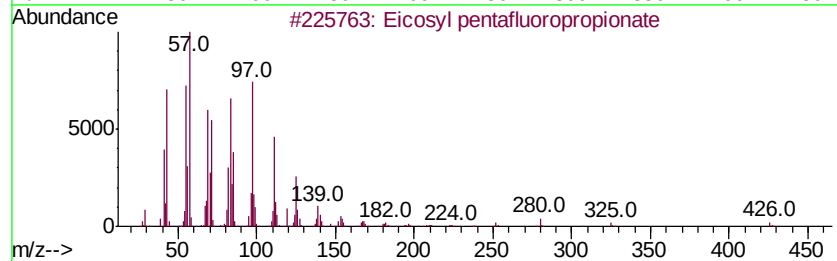
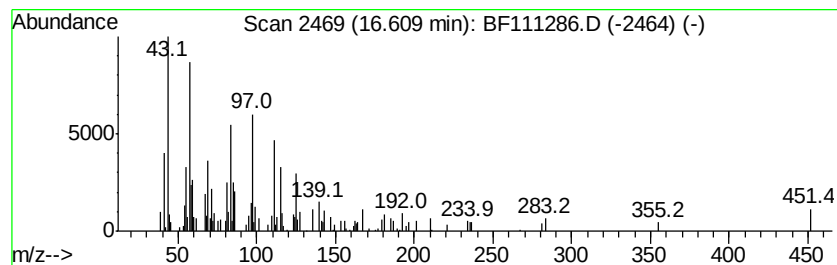
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Eicosyl pentafluoropropionate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.61	3.27 ng	154130	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	50
2		Heneicosyl trifluoroacetate	408	C23H43F3O2	1000351-76-5	50
3		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	46
4		Tetracosyl heptafluorobutyrte	550	C28H49F7O2	1000351-83-7	46
5		Tricosyl heptafluorobutyrate	536	C27H47F7O2	1000351-83-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111286.D
Acq On : 11 Dec 2018 19:36
Operator : JU/SJ
Sample : J6157-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMPOSITE-2-DISSOLVED

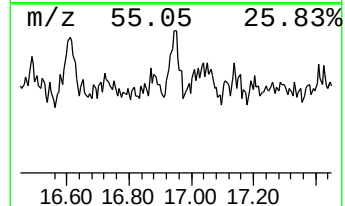
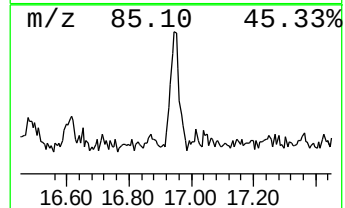
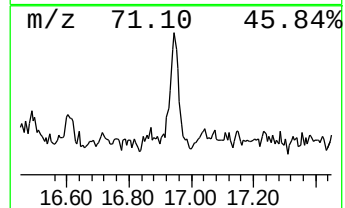
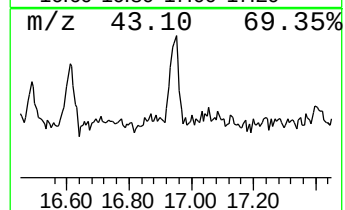
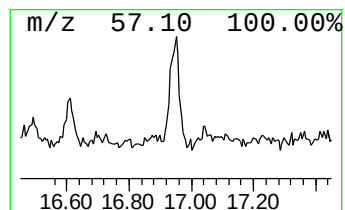
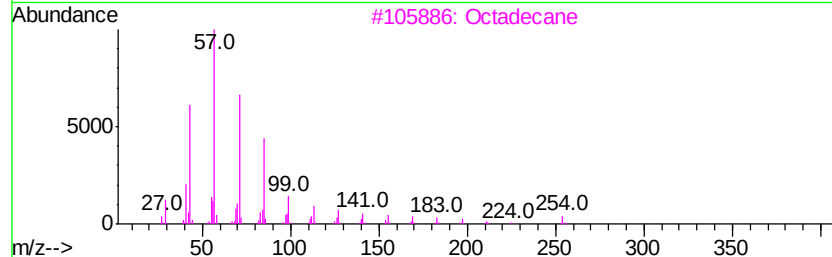
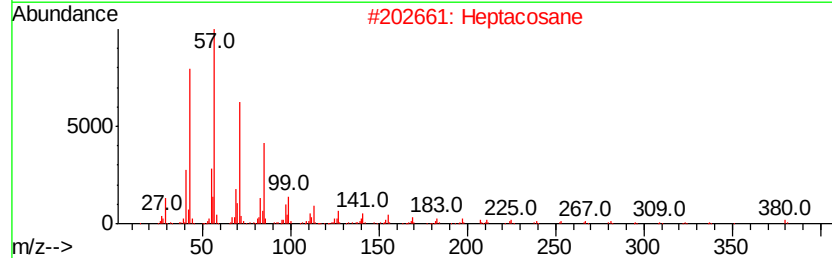
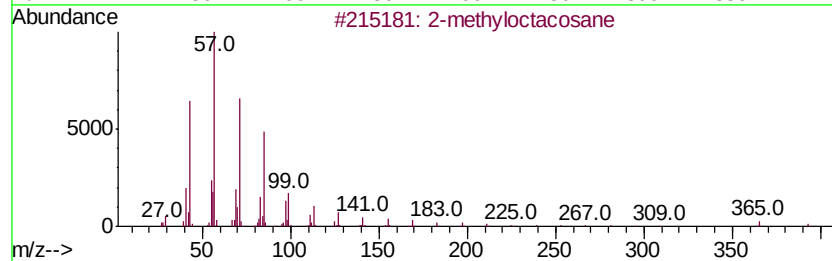
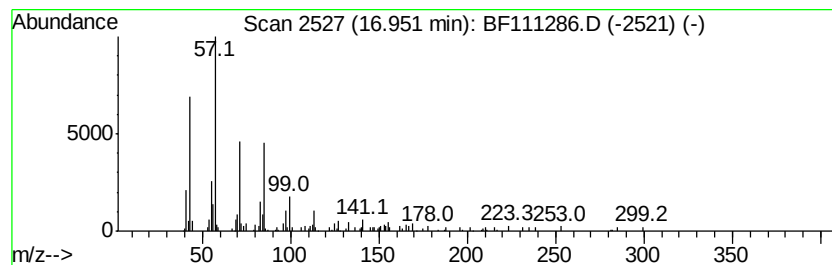
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2-methyloctacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.95	3.16 ng	148750	Perylene-d12	15.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-methyloctacosane	408	C29H60	1000376-72-8	86
2		Heptacosane	380	C27H56	000593-49-7	86
3		Octadecane	254	C18H38	000593-45-3	80
4		Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	64
5		Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
Data File : BF111286.D
Acq On : 11 Dec 2018 19:36
Operator : JU/SJ
Sample : J6157-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

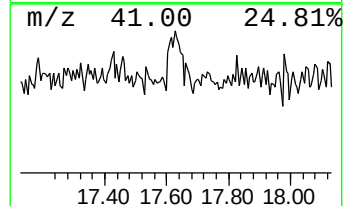
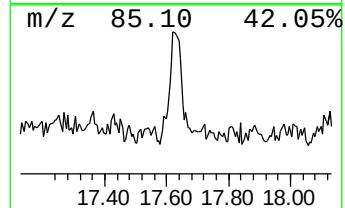
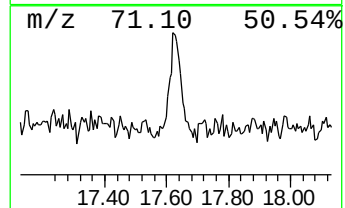
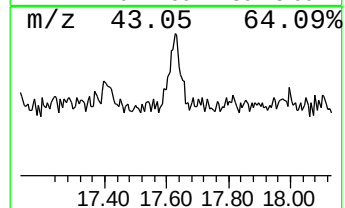
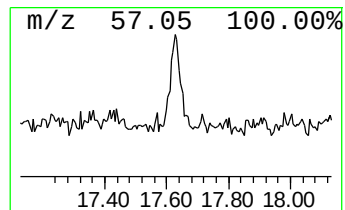
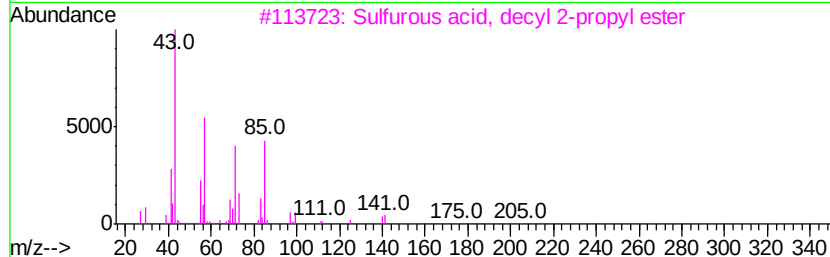
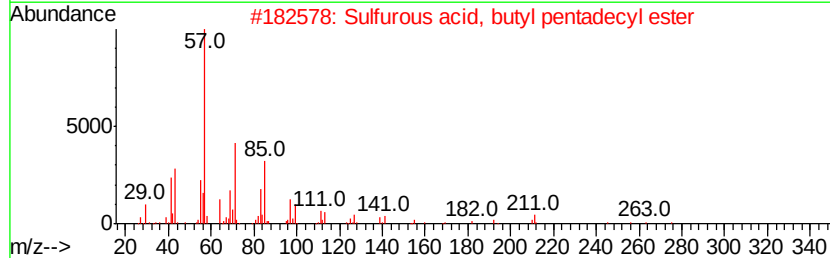
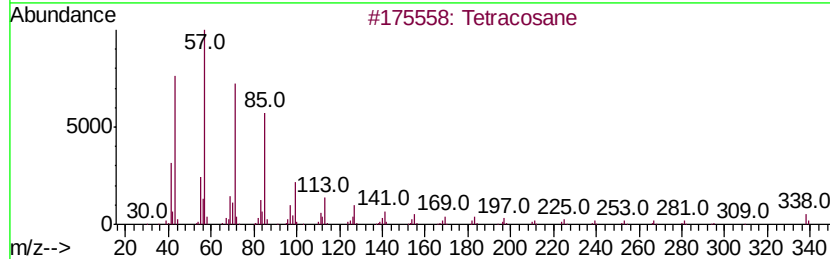
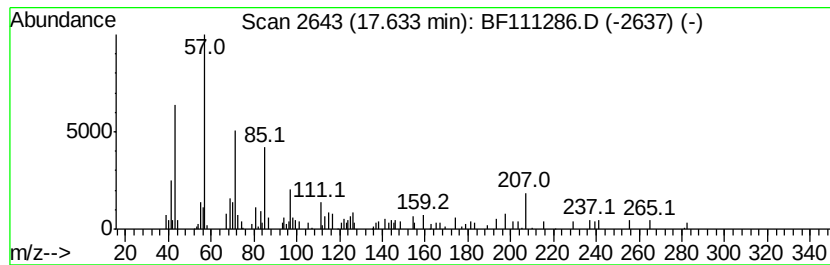
Instrument :
BNA_F
ClientSampleId :
COMPOSITE-2-DISSOLVED

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetracosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.63	2.85 ng	134342	Perylene-d12	15.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetracosane	338 C24H50	000646-31-1 50
2		Sulfurous acid, butyl pentadecyl...	348 C19H40O3S	1000309-18-2 50
3		Sulfurous acid, decyl 2-propyl e...	264 C13H28O3S	1000309-12-1 50
4		Octane, 3,4,5,6-tetramethyl-	170 C12H26	062185-21-1 50
5		Tridecane, 1-iodo-	310 C13H27I	035599-77-0 50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121118\
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 Operator : JU/SJ
 Sample : J6157-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMPOSITE-2-DISSOLVED

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120818.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
1,3-Dioxolane, 2,...	2.83	5.4	ng	390186	1	6.80	1448290	20.0
2-Pentanone, 4-hy...	4.99	4.4	ng	317497	1	6.80	1448290	20.0
unknown6.56	6.56	56.4	ng	4086010	1	6.80	1448290	20.0
n-Hexadecanoic acid	11.86	20.7	ng	2341640	4	11.32	2259630	20.0
Octadecanoic acid	12.64	13.9	ng	1379570	5	13.96	1991130	20.0
Pentafluoropropio...	13.81	11.2	ng	1119060	5	13.96	1991130	20.0
Octacosane	14.44	3.2	ng	314498	5	13.96	1991130	20.0
Docosane	15.08	4.0	ng	189804	6	15.40	941594	20.0
Heptadecane, 9-oc...	15.87	4.8	ng	225203	6	15.40	941594	20.0
10-Heneicosene (c,t)	15.93	2.1	ng	99548	6	15.40	941594	20.0
Tridecane, 1-iodo-	16.37	4.8	ng	226288	6	15.40	941594	20.0
Eicosyl pentaflu...	16.61	3.3	ng	154130	6	15.40	941594	20.0
2-methyloctacosane	16.95	3.2	ng	148750	6	15.40	941594	20.0
Tetracosane	17.63	2.9	ng	134342	6	15.40	941594	20.0