

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
 Data File : BF119426.D
 Acq On : 18 Mar 2020 20:28
 Operator : CG/JU
 Sample : L1942-21
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-5-SA

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	2.151	7	11	24	rVB	302681	430714	5.17%	0.656%
2	5.075	501	508	515	rVB	148452	196671	2.36%	0.300%
3	5.469	569	575	578	rBV	7126822	7622011	91.46%	11.616%
4	6.475	740	746	749	rBV	6876456	7639199	91.67%	11.642%
5	6.851	807	810	814	rVB	2492847	2161411	25.94%	3.294%
6	7.010	833	837	841	rVB	6654395	5985827	71.83%	9.122%
7	7.416	900	906	909	rBV	5171573	5019818	60.24%	7.650%
8	8.139	1024	1029	1032	rBV	3378798	2841648	34.10%	4.331%
9	9.222	1207	1213	1216	rBV	8822817	8333323	100.00%	12.700%
10	9.610	1275	1279	1282	rBV	892218	661138	7.93%	1.008%
11	9.898	1323	1328	1331	rBV	3694023	3063297	36.76%	4.669%
12	10.192	1374	1378	1386	rVB	160853	148475	1.78%	0.226%
13	10.686	1456	1462	1465	rBV	5553654	5010275	60.12%	7.636%
14	11.386	1577	1581	1584	rBV	3661933	3013536	36.16%	4.593%
15	11.910	1664	1670	1675	rBV	242392	245022	2.94%	0.373%
16	12.698	1801	1804	1810	rBV2	109404	113466	1.36%	0.173%
17	12.974	1847	1851	1855	rBV	7334927	7041257	84.50%	10.731%
18	13.874	2000	2004	2010	rBV	513924	581405	6.98%	0.886%
19	14.021	2025	2029	2033	rBV	2196178	2056962	24.68%	3.135%
20	14.198	2056	2059	2064	rVB	168650	152634	1.83%	0.233%
21	14.504	2108	2111	2115	rBV	271343	244276	2.93%	0.372%
22	14.815	2160	2164	2169	rVB	284275	290118	3.48%	0.442%
23	15.162	2218	2223	2228	rBV	255968	301702	3.62%	0.460%
24	15.492	2274	2279	2284	rBV2	1469617	1768666	21.22%	2.695%
25	15.545	2284	2288	2296	rVB	225937	292076	3.50%	0.445%
26	15.992	2358	2364	2368	rBV	150738	241821	2.90%	0.369%
27	16.503	2446	2451	2457	rBV2	85098	159365	1.91%	0.243%

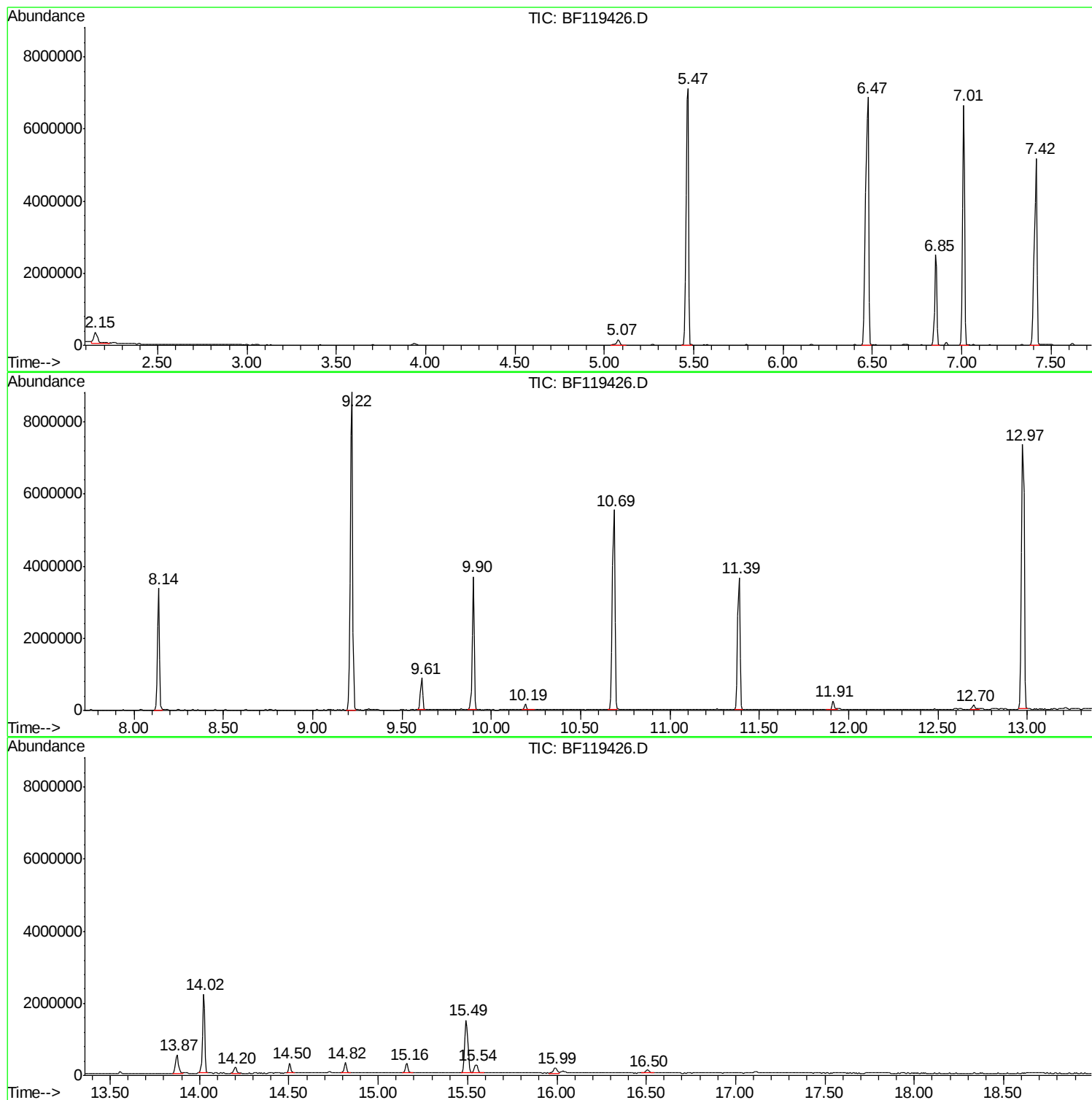
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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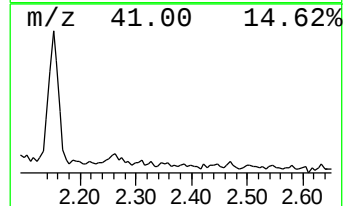
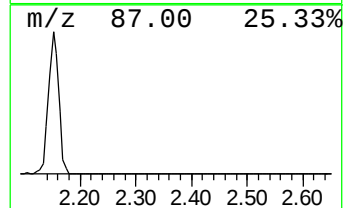
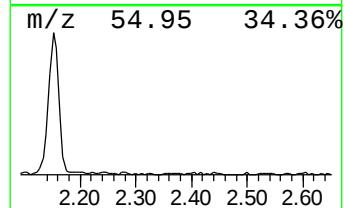
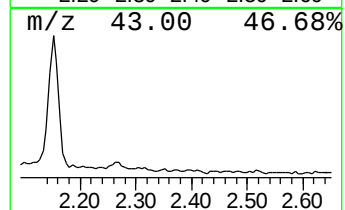
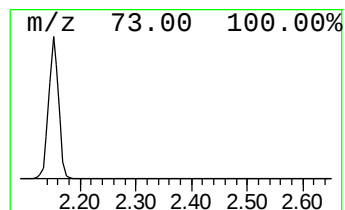
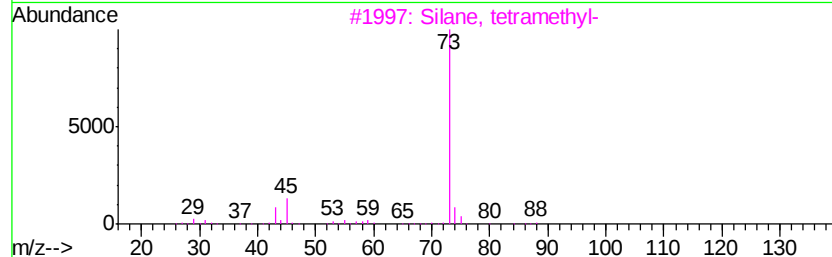
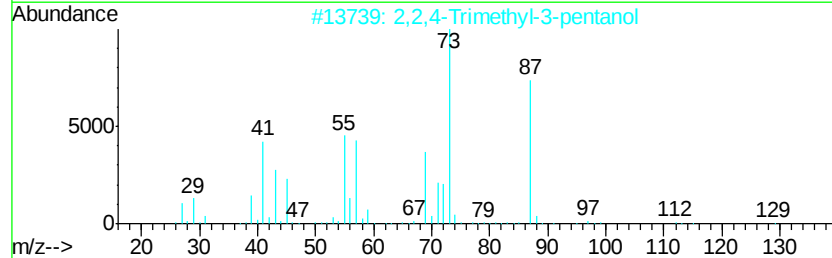
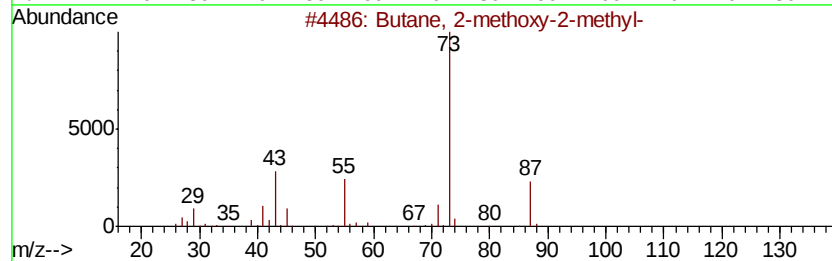
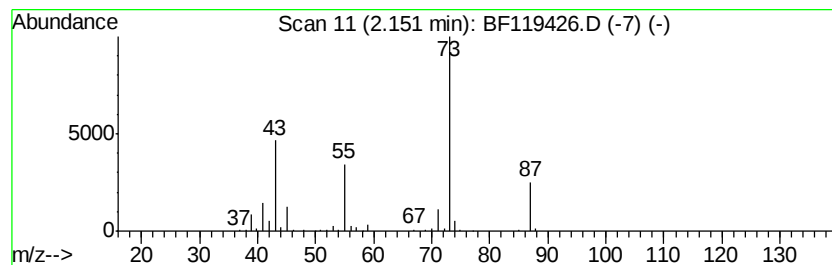
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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	3.99 ng	430714	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		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
4		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	23
5		3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	12



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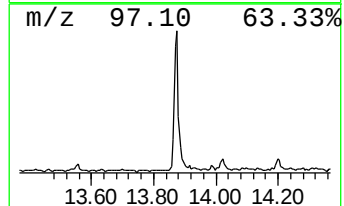
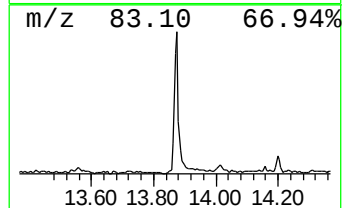
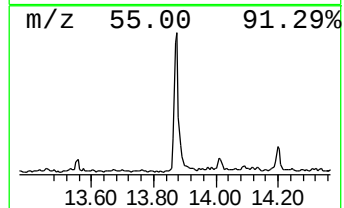
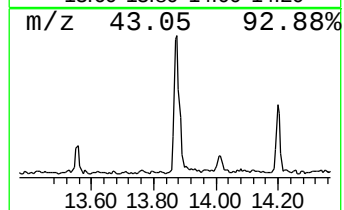
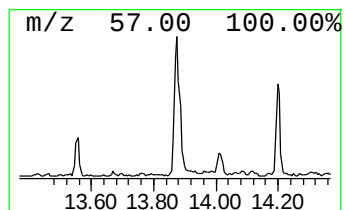
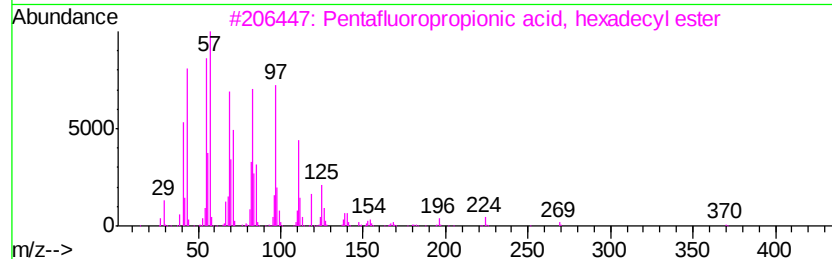
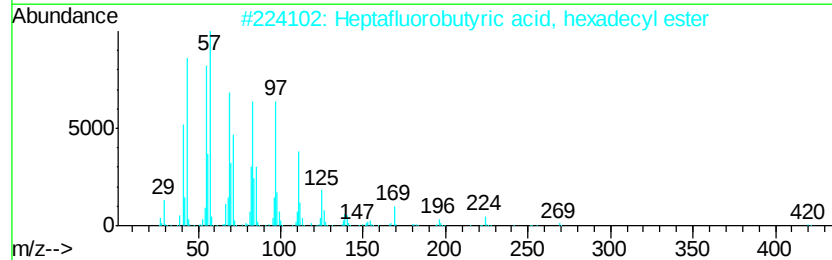
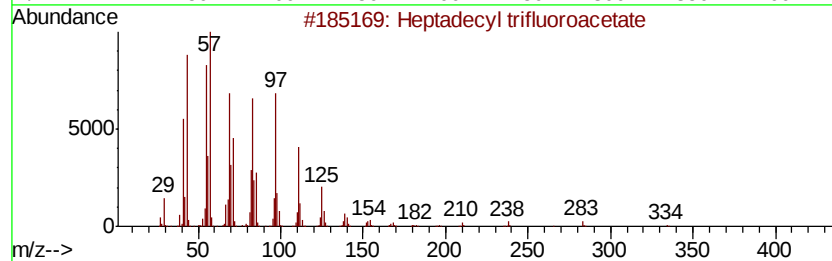
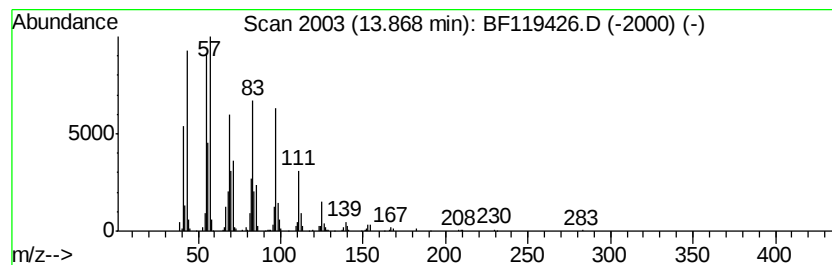
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Heptadecyl trifluoroacetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.87	5.65 ng	581405	Chrysene-d12	14.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	94
2		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	94
3		Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
4		1-Hexadecanol	242	C16H34O	036653-82-4	91
5		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



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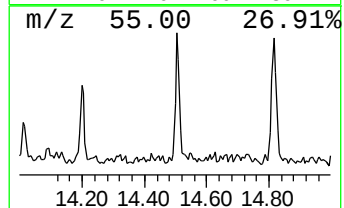
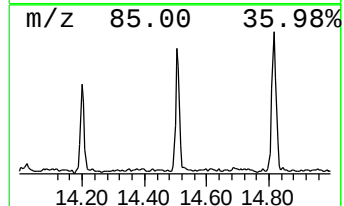
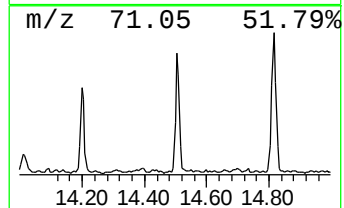
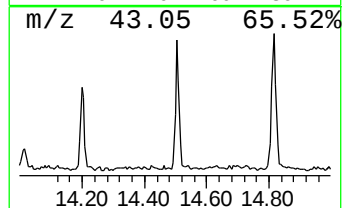
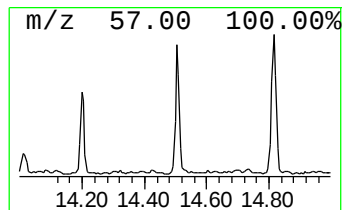
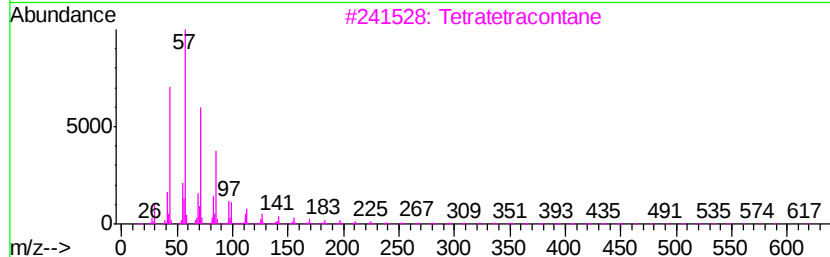
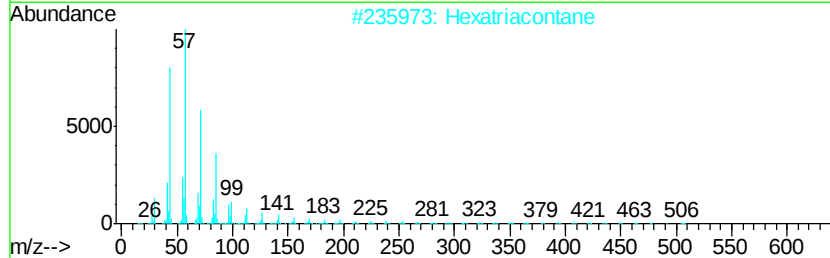
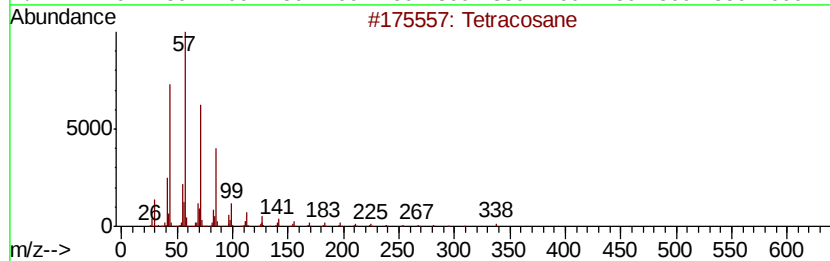
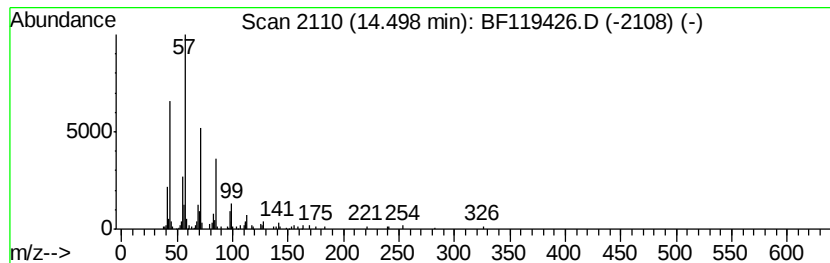
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 TIC Integration Parameters: LSCINT.P

 Peak Number 4 Tetracosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	2.38 ng	244276	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	91
2		Hexatriacontane	507	C36H74	000630-06-8	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		Hexacosane	366	C26H54	000630-01-3	90
5		Octacosane	394	C28H58	000630-02-4	90



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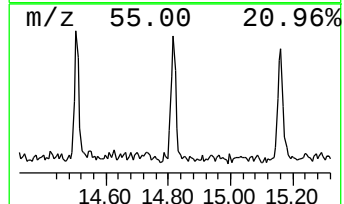
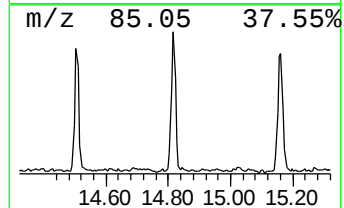
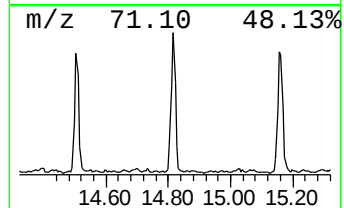
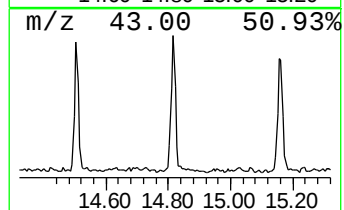
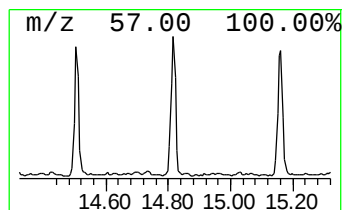
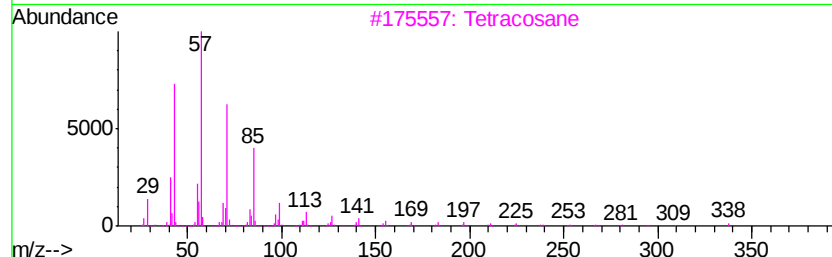
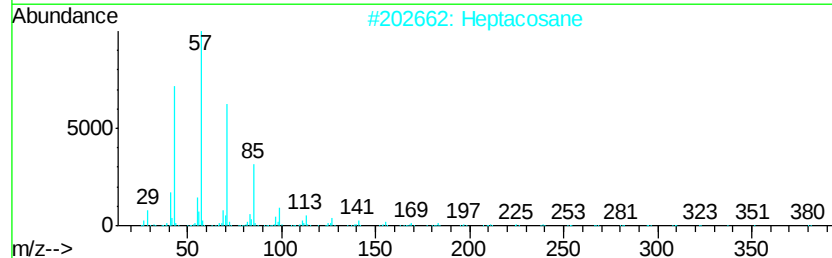
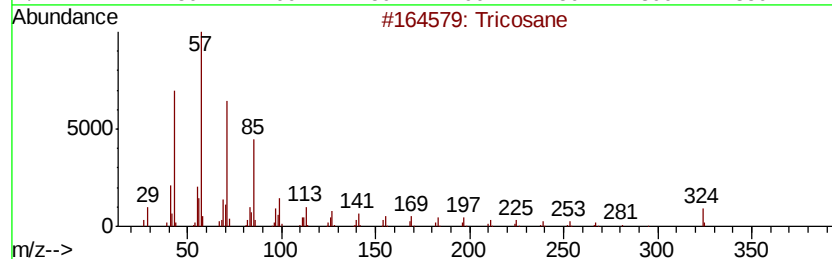
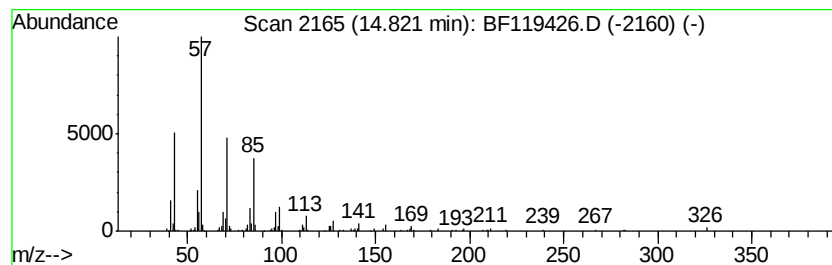
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tricosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	3.28 ng	290118	Perylene-d12	15.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane		324	C23H48	000638-67-5	94
2	Heptacosane		380	C27H56	000593-49-7	91
3	Tetracosane		338	C24H50	000646-31-1	91
4	Hexadecane		226	C16H34	000544-76-3	91
5	Heneicosane		296	C21H44	000629-94-7	90



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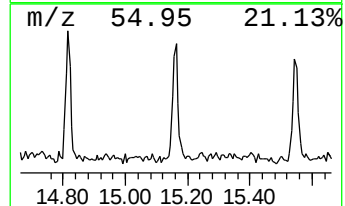
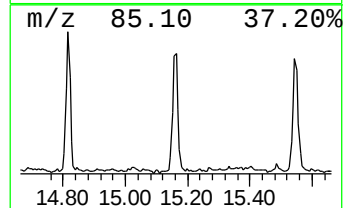
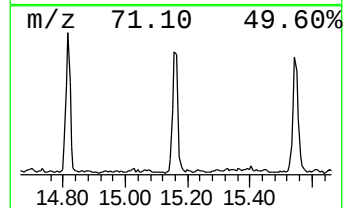
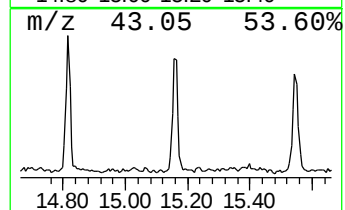
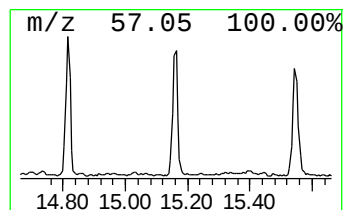
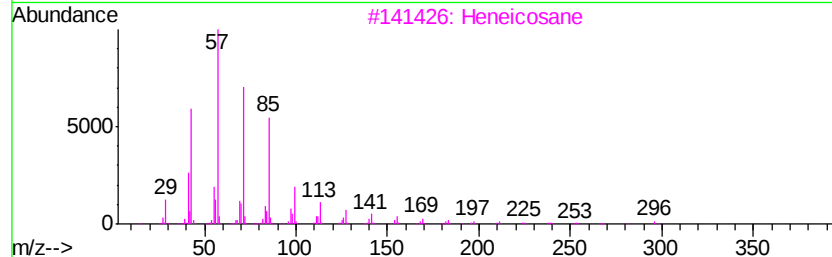
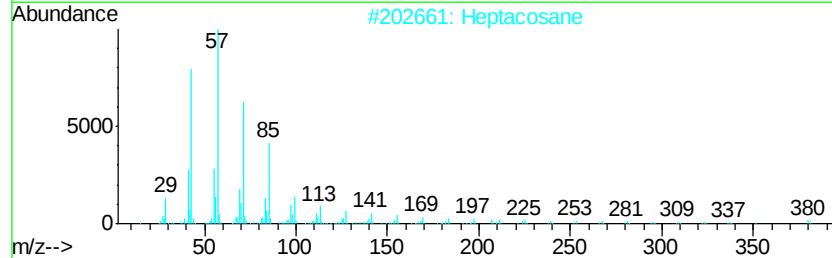
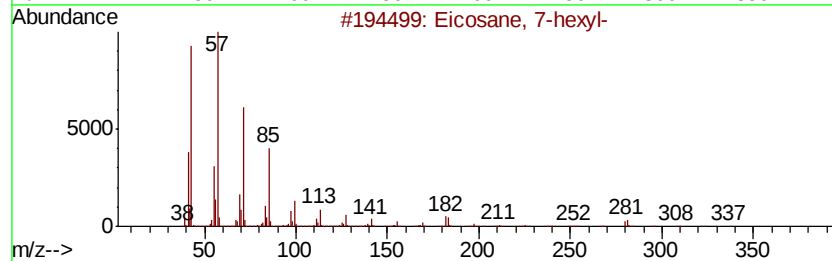
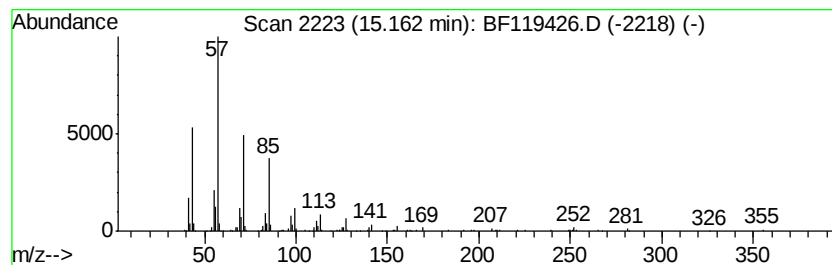
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Eicosane, 7-hexyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.16	3.41 ng	301702	Perylene-d12		15.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 7-hexyl-			366	C26H54	055333-99-8	90
2	Heptacosane			380	C27H56	000593-49-7	87
3	Heneicosane			296	C21H44	000629-94-7	87
4	Hexacosane			366	C26H54	000630-01-3	87
5	Pentadecane			212	C15H32	000629-62-9	86



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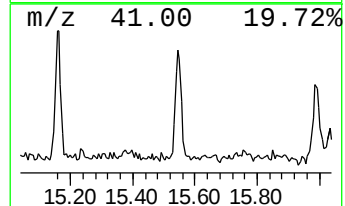
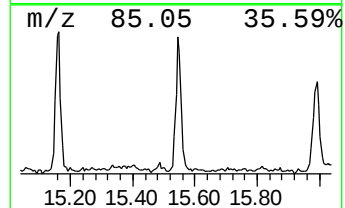
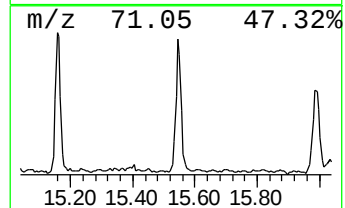
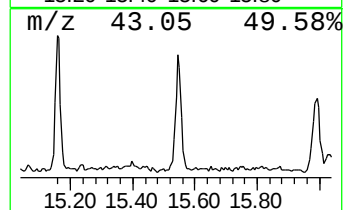
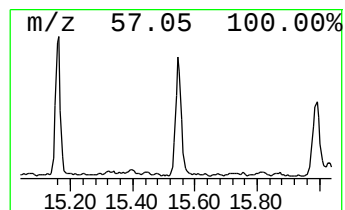
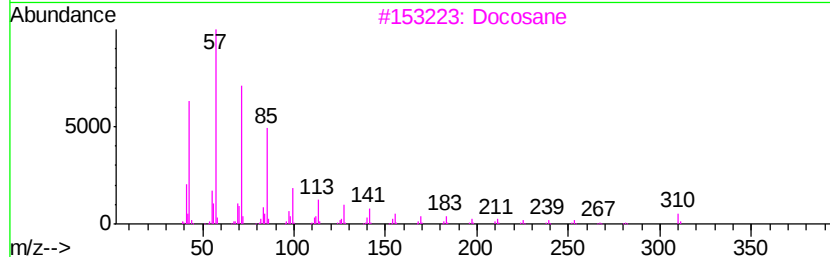
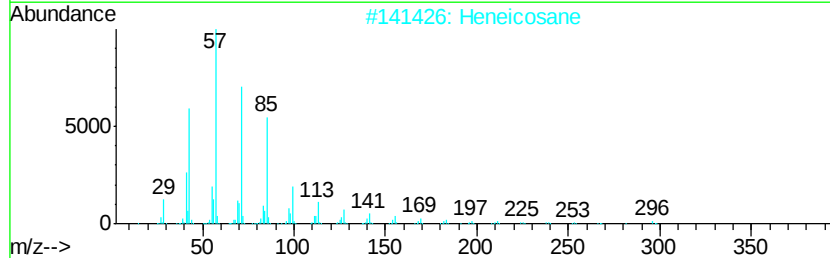
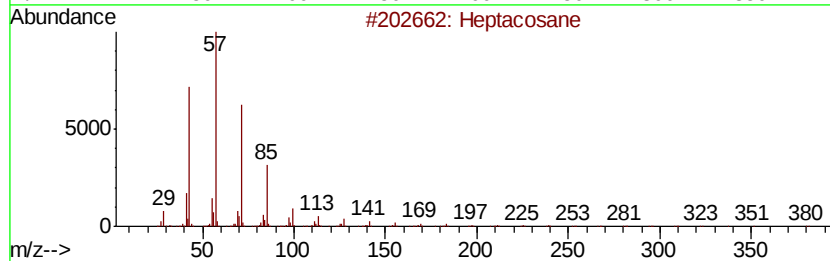
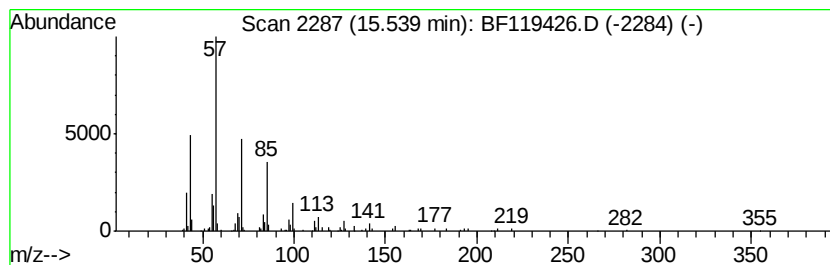
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	3.30 ng	292076	Perylene-d12	15.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Docosane	310	C22H46	000629-97-0	90
4		Heptadecane	240	C17H36	000629-78-7	90
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031820\
Data File : BF119426.D
Acq On : 18 Mar 2020 20:28
Operator : CG/JU
Sample : L1942-21
Misc :
ALS Vial : 11 Sample Multiplier: 1

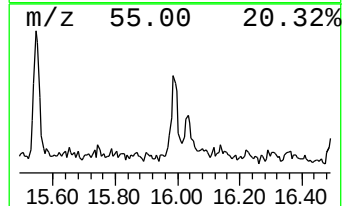
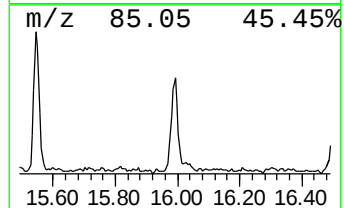
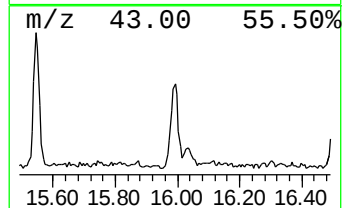
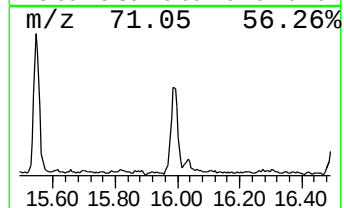
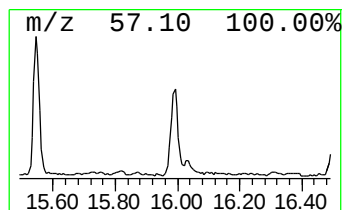
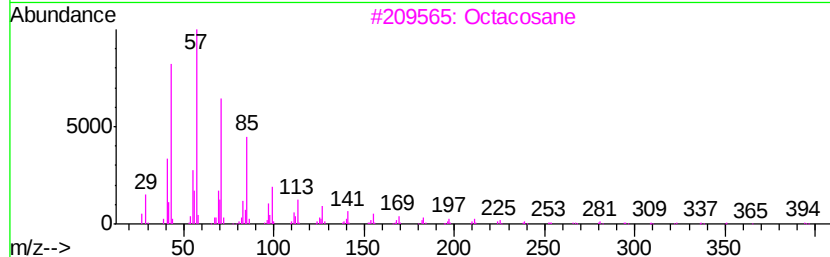
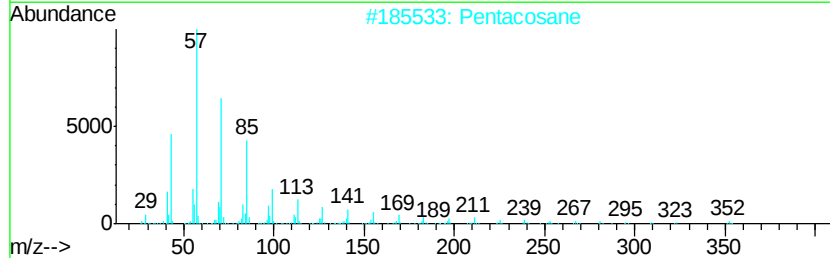
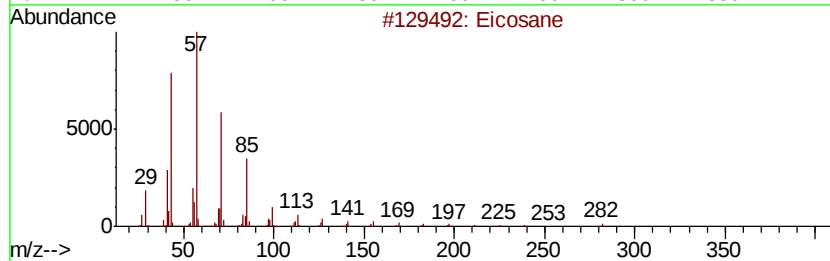
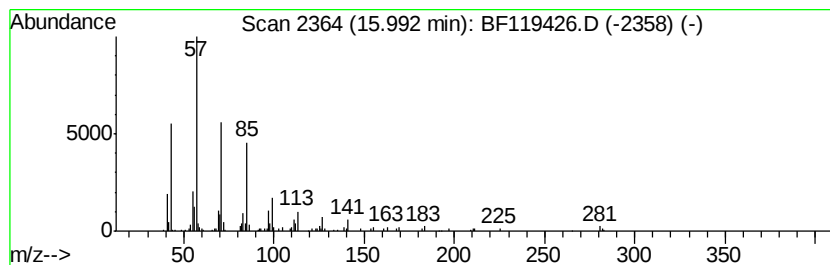
Instrument :
BNA_F
ClientSampleId :
TP-5-SA

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

Peak Number 8 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	2.73 ng	241821	Perylene-d12	15.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 93
2	Pentacosane		352 C25H52	000629-99-2 91
3	Octacosane		394 C28H58	000630-02-4 90
4	2-Bromotetradecane		276 C14H29Br	074036-95-6 87
5	Octadecane, 1-iodo-		380 C18H37I	000629-93-6 86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031820\
Data File : BF119426.D
Acq On : 18 Mar 2020 20:28
Operator : CG/JU
Sample : L1942-21
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-5-SA

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...	2.15	4.0	ng	430714	1	6.85	2161410	20.0
Heptadecyl triflu...	13.87	5.7	ng	581405	5	14.02	2056960	20.0
Tetracosane	14.50	2.4	ng	244276	5	14.02	2056960	20.0
Tricosane	14.82	3.3	ng	290118	6	15.49	1768670	20.0
Eicosane, 7-hexyl-	15.16	3.4	ng	301702	6	15.49	1768670	20.0
Heptacosane	15.54	3.3	ng	292076	6	15.49	1768670	20.0
Eicosane	15.99	2.7	ng	241821	6	15.49	1768670	20.0