

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF123119\
 Data File : BF118402.D
 Acq On : 31 Dec 2019 13:44
 Operator : JU
 Sample : K6463-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-7-(4-6)

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-BF122419.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.057	501	505	518	rVB	280155	354706	10.13%	1.042%
2	5.475	572	576	597	rBV	2108591	2188570	62.52%	6.430%
3	6.492	744	749	767	rBV	2157251	2357050	67.34%	6.925%
4	6.622	767	771	775	rBV	2711157	2478423	70.81%	7.281%
5	6.851	806	810	818	rBV	917256	719873	20.57%	2.115%
6	7.010	833	837	840	rBV	2138394	1927269	55.06%	5.662%
7	7.416	902	906	920	rBV	1488957	1452842	41.51%	4.268%
8	8.139	1025	1029	1044	rBV	1016435	942053	26.91%	2.768%
9	9.216	1208	1212	1227	rBV	3331796	3069453	87.69%	9.017%
10	9.616	1276	1280	1291	rBV	150369	161955	4.63%	0.476%
11	9.898	1324	1328	1339	rBV	1165887	1152978	32.94%	3.387%
12	10.698	1460	1464	1477	rBV	2158789	2122140	60.63%	6.234%
13	11.392	1579	1582	1591	rBV	1231273	1243688	35.53%	3.654%
14	11.916	1666	1671	1679	rBV2	349316	408836	11.68%	1.201%
15	12.621	1786	1791	1797	rBV4	32996	61216	1.75%	0.180%
16	12.704	1800	1805	1818	rBV	271636	372953	10.65%	1.096%
17	12.986	1849	1853	1856	rBV	4009098	3500334	100.00%	10.283%
18	13.204	1886	1890	1895	rBV2	39001	48037	1.37%	0.141%
19	13.545	1944	1948	1951	rBV2	59300	66327	1.89%	0.195%
20	13.880	2001	2005	2017	rBV	404681	555243	15.86%	1.631%
21	14.051	2030	2034	2043	rBV	1116248	1075927	30.74%	3.161%
22	14.198	2054	2059	2070	rVB	399350	424495	12.13%	1.247%
23	14.504	2106	2111	2114	rBV	757687	768240	21.95%	2.257%
24	14.662	2135	2138	2142	rBV	191544	197900	5.65%	0.581%
25	14.809	2158	2163	2173	rVB	1000213	1152759	32.93%	3.387%
26	15.157	2216	2222	2229	rBV	1008467	1243522	35.53%	3.653%
27	15.545	2281	2288	2306	rVB2	1397095	2213118	63.23%	6.502%
28	15.809	2330	2333	2337	rVB4	31062	42466	1.21%	0.125%
29	15.986	2355	2363	2370	rBV	595273	1015357	29.01%	2.983%
30	16.498	2443	2450	2456	rBV	299340	525538	15.01%	1.544%
31	17.098	2546	2552	2560	rVB	90074	195829	5.59%	0.575%

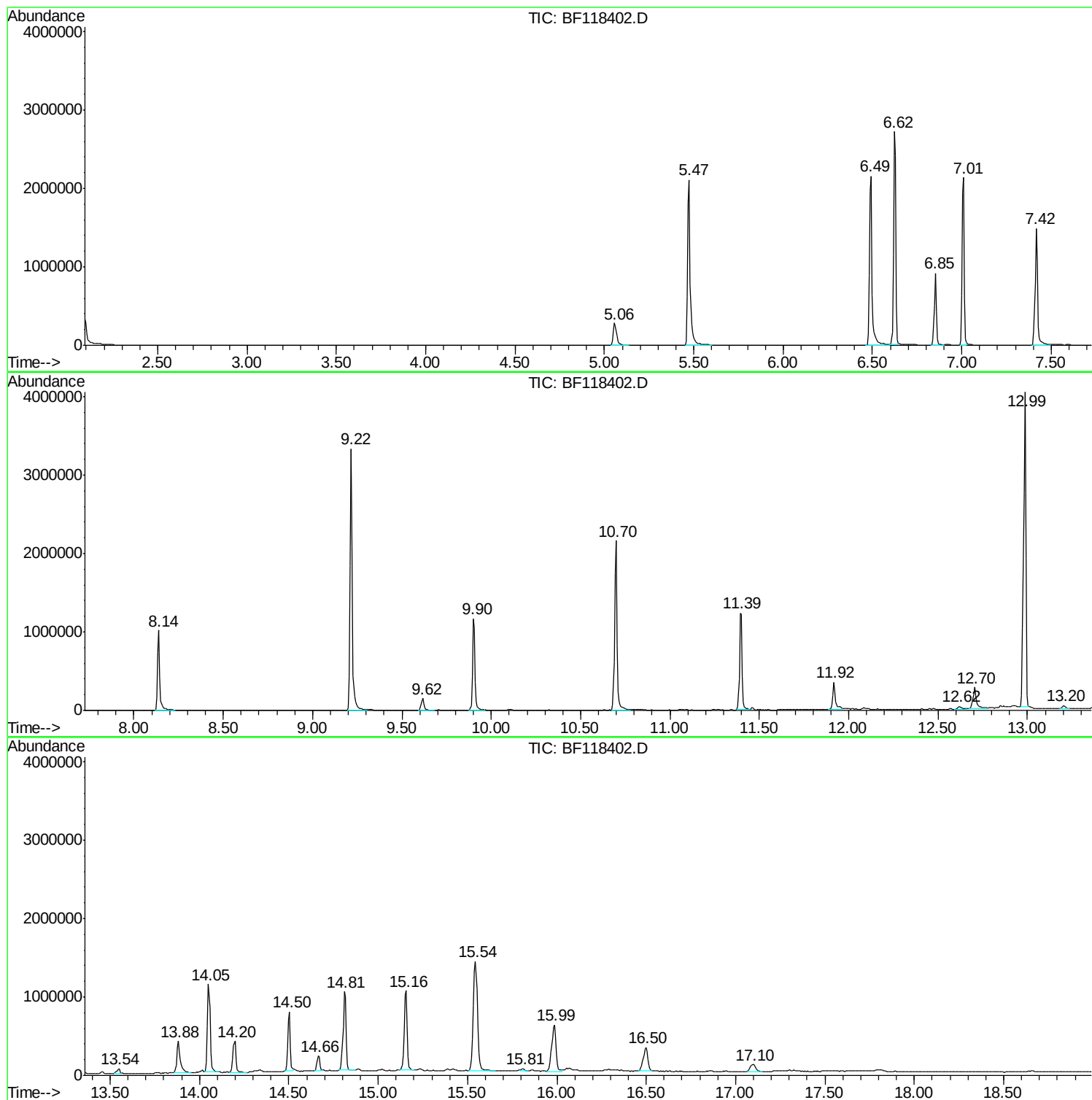
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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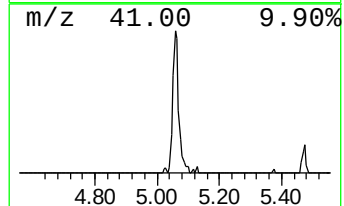
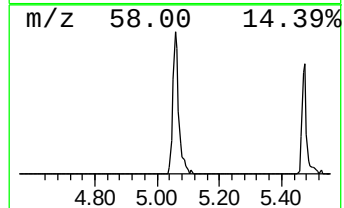
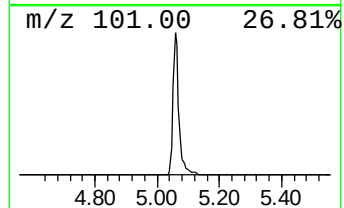
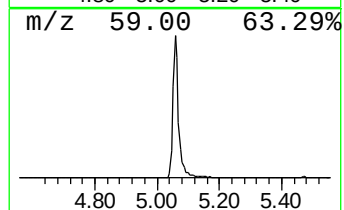
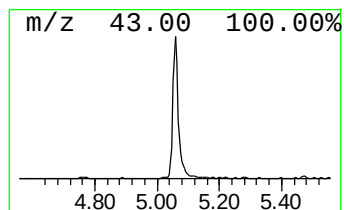
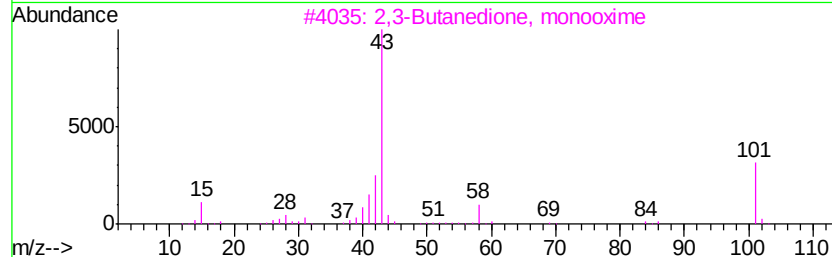
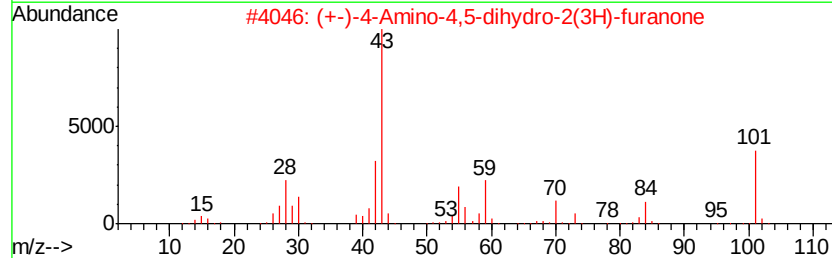
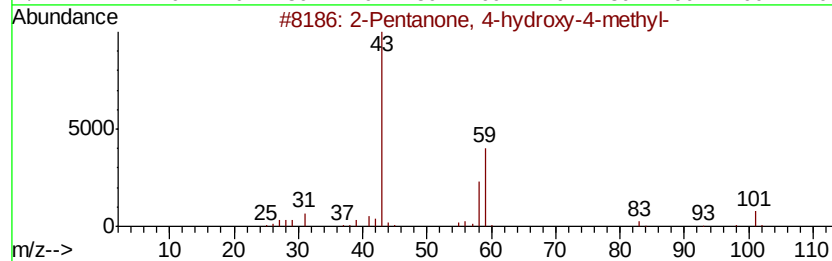
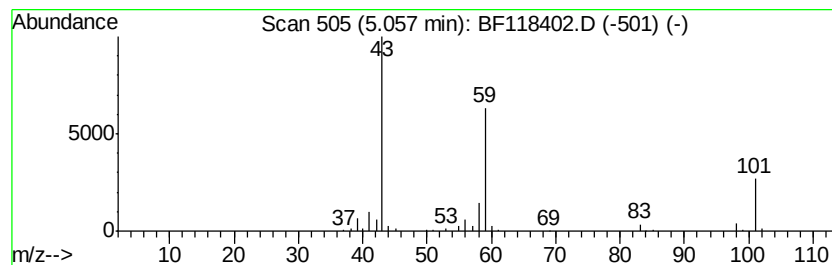
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	9.85 ng	354706	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	40
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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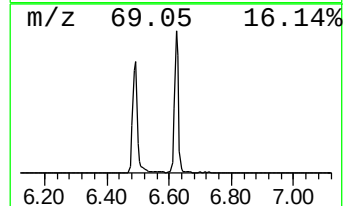
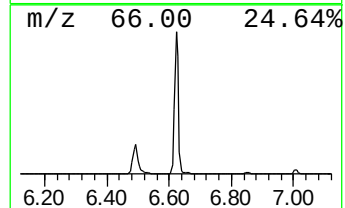
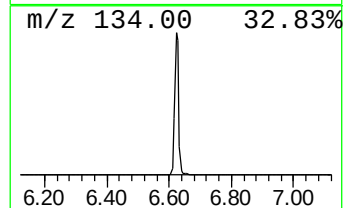
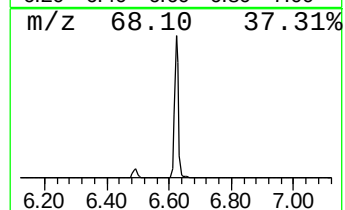
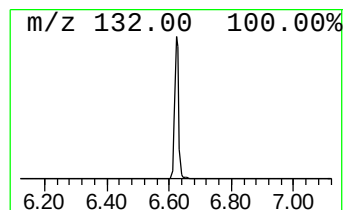
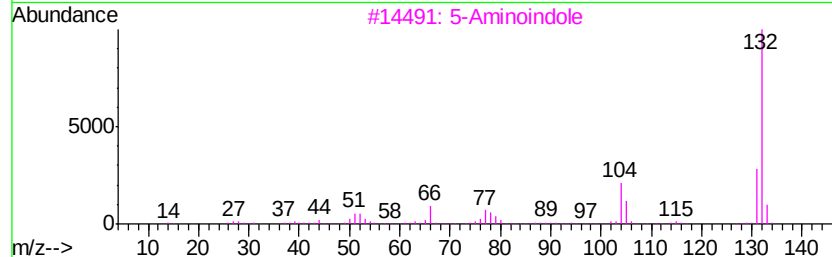
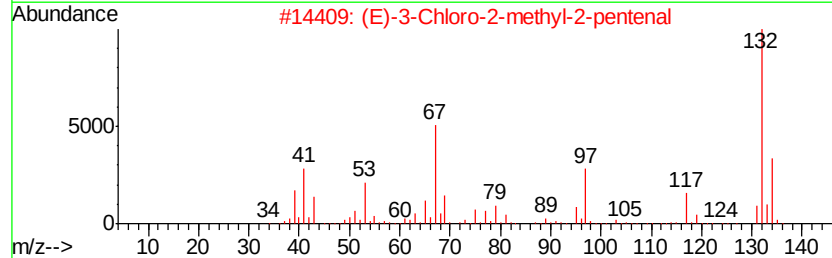
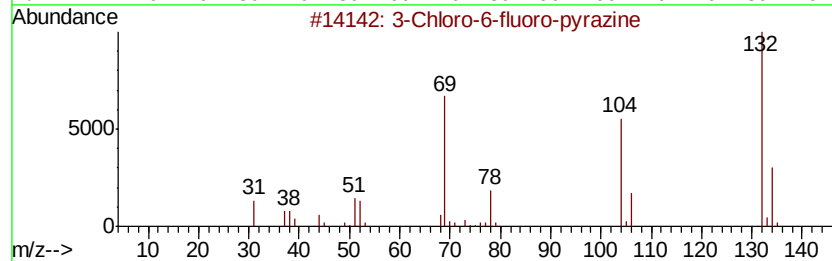
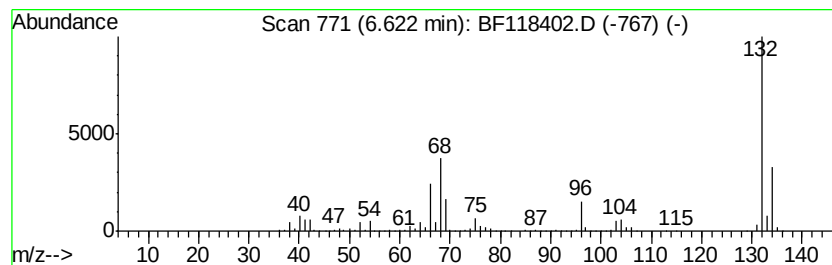
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	68.86 ng	2478420	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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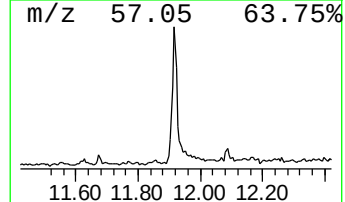
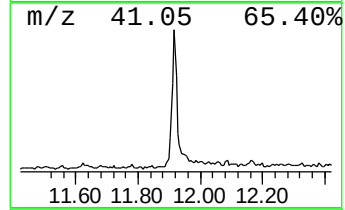
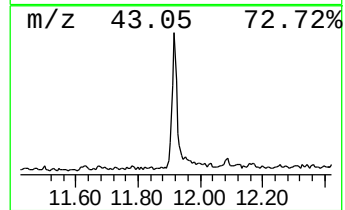
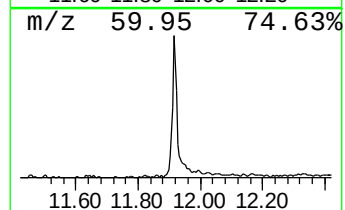
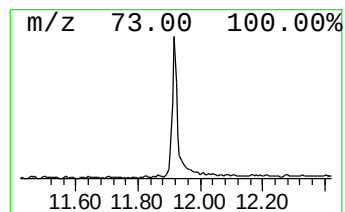
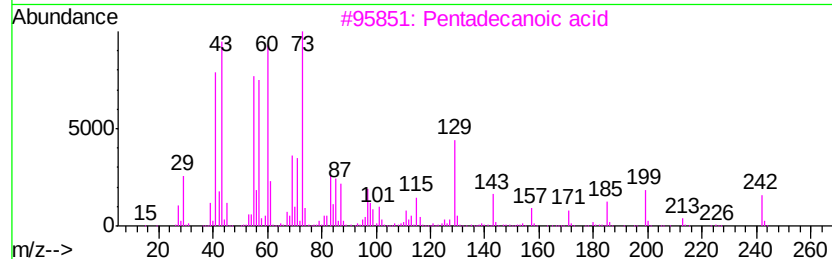
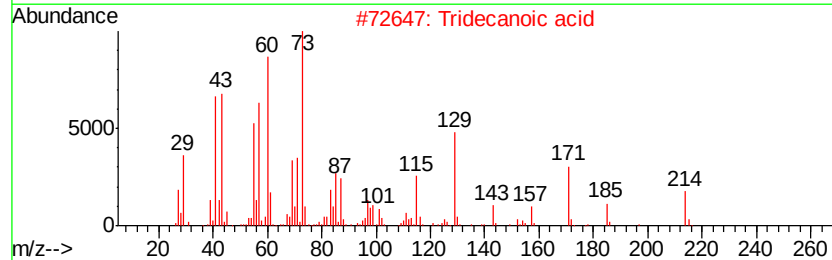
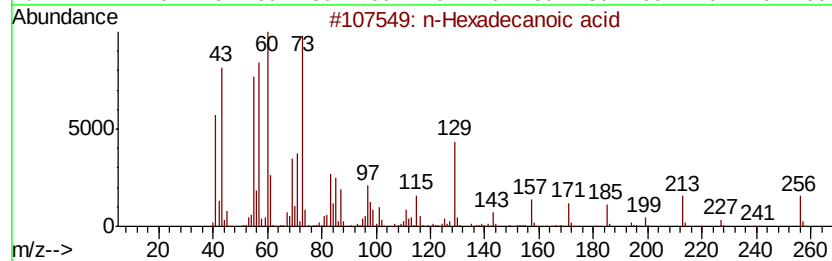
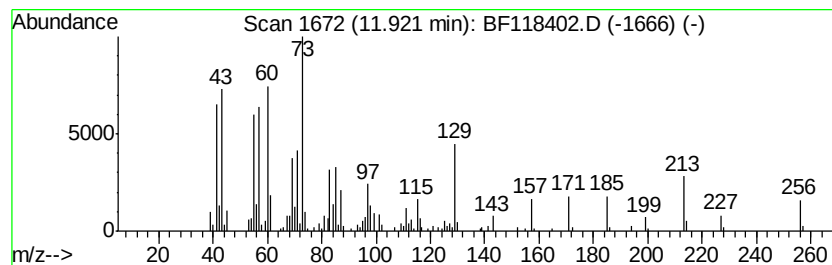
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.92	6.57 ng	408836	Phenanthrene-d10		11.40
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	94
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
5	Undecanoic acid	186	C11H22O2	000112-37-8	50



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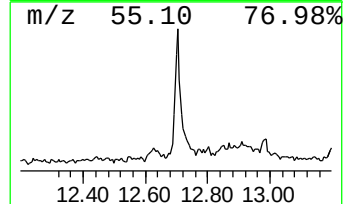
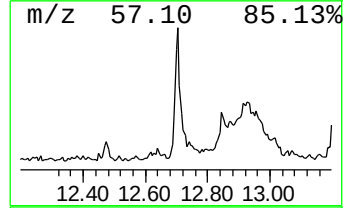
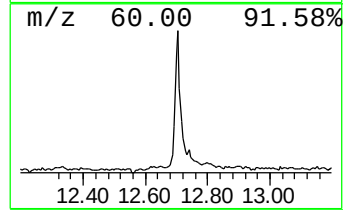
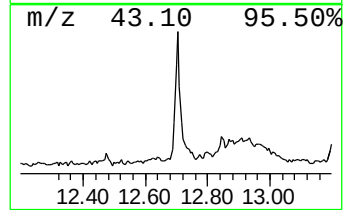
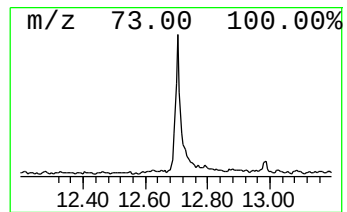
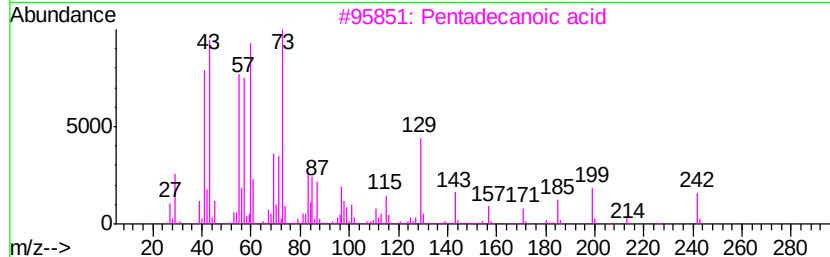
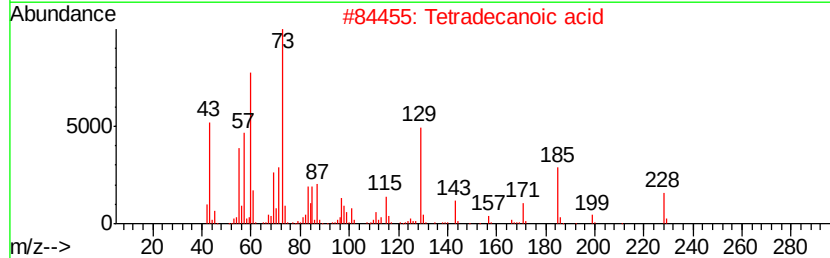
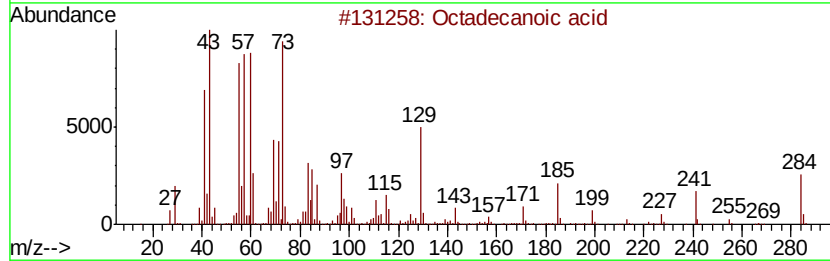
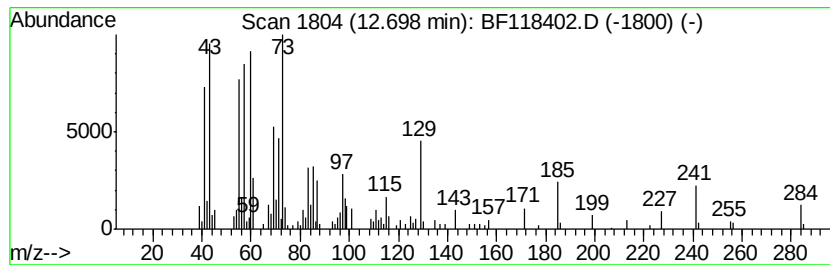
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Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.70	6.00 ng	372953	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	62
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
4	n-Decanoic acid	172	C10H20O2	000334-48-5	55
5	Dodecanoic acid	200	C12H24O2	000143-07-7	50



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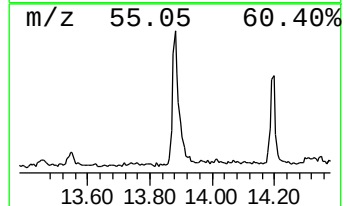
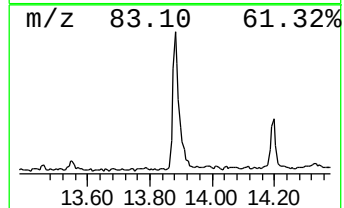
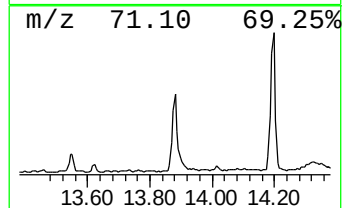
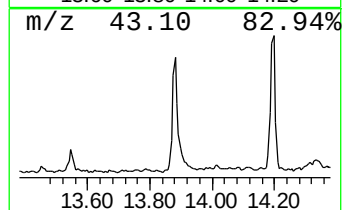
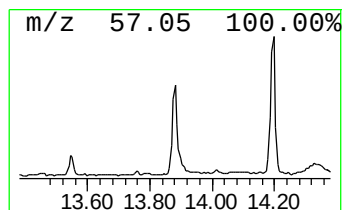
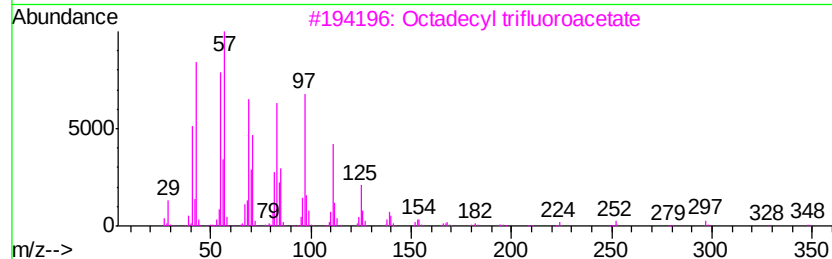
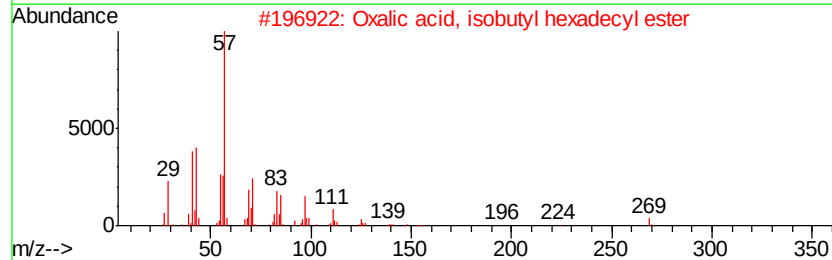
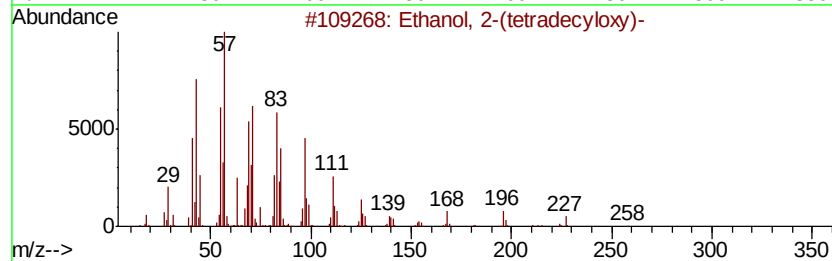
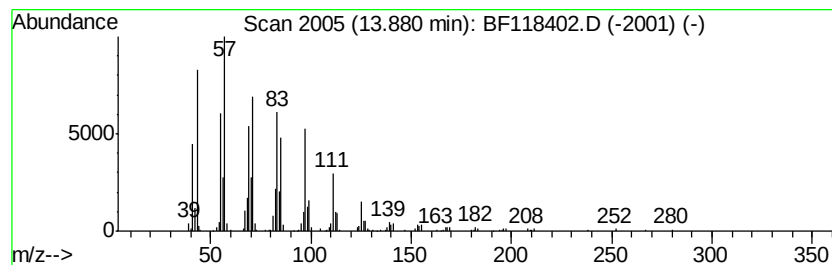
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Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.88	10.32 ng	555243	Chrysene-d12	14.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	76
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	76
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	76



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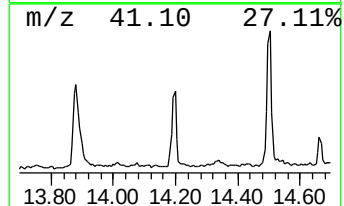
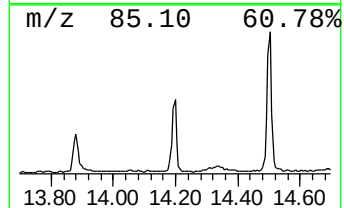
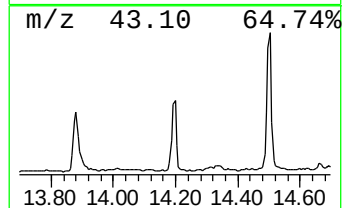
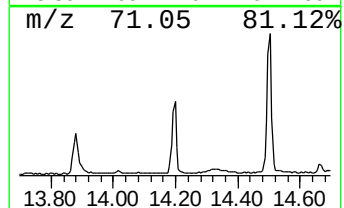
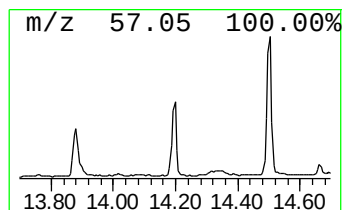
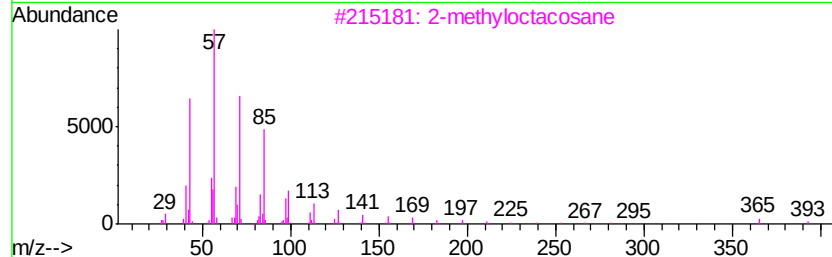
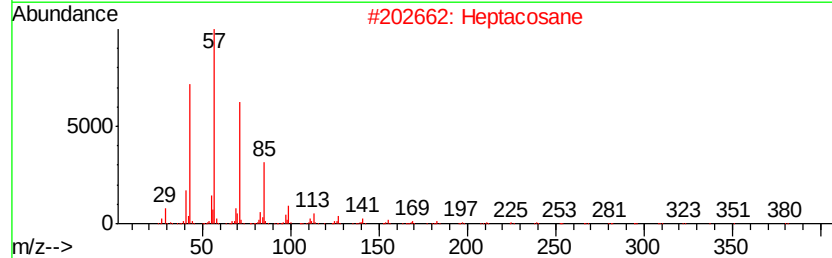
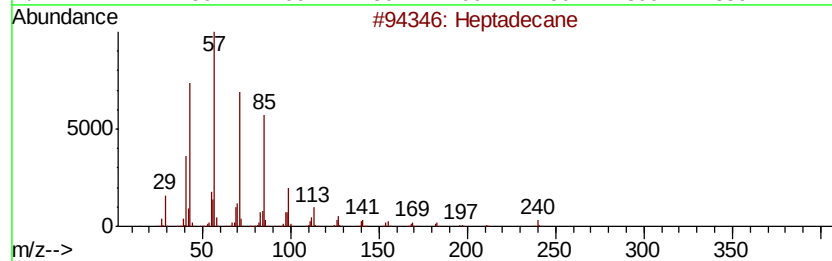
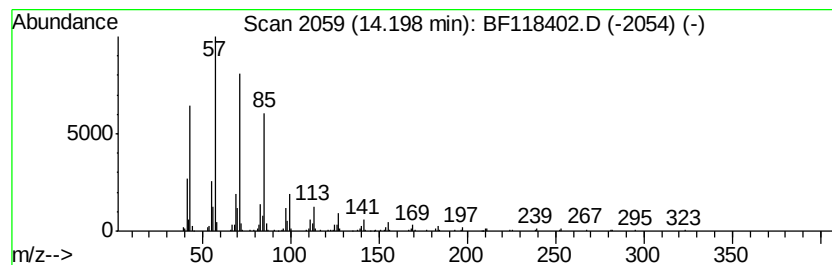
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	7.89 ng	424495	Chrysene-d12	14.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Heptacosane		380	C27H56	000593-49-7	91
3	2-methyloctacosane		408	C29H60	1000376-72-8	91
4	Heneicosane		296	C21H44	000629-94-7	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
 Data File : BF118402.D
 Acq On : 31 Dec 2019 13:44
 Operator : JU
 Sample : K6463-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

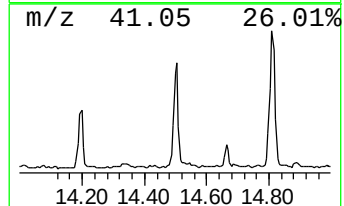
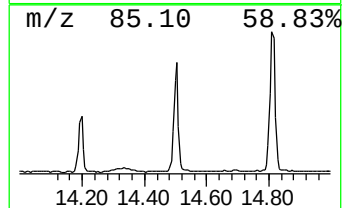
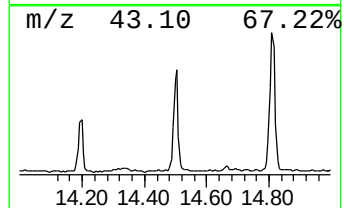
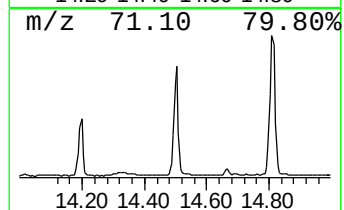
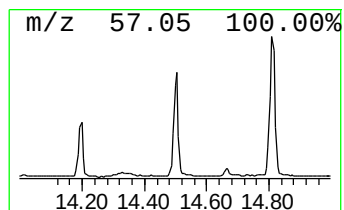
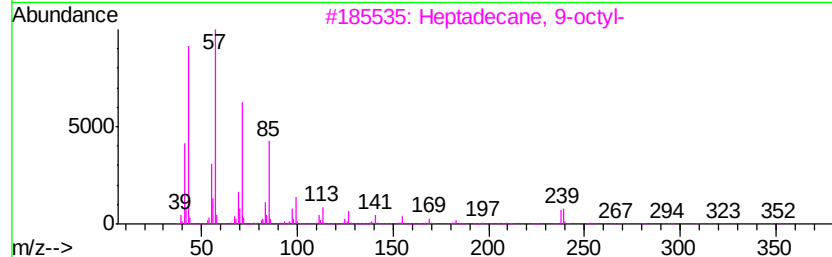
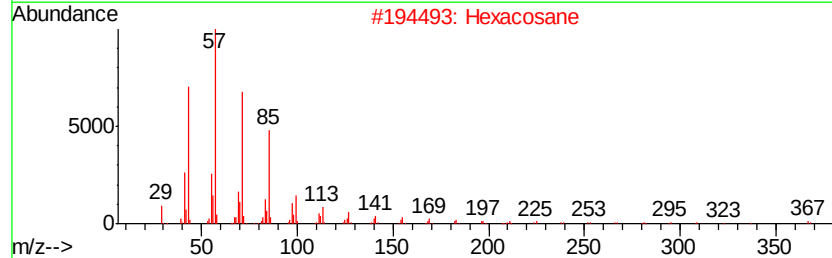
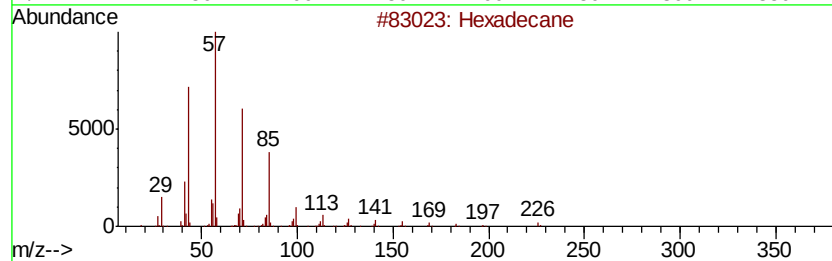
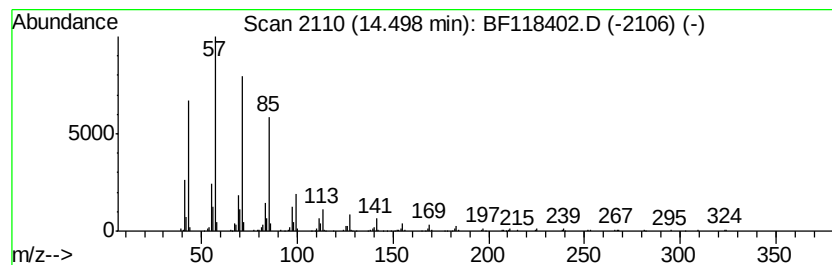
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 ClientSampleId :
 GP-7-(4-6)

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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 Hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	14.28 ng	768240	Chrysene-d12	14.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	95
2	Hexacosane	366 C26H54	000630-01-3	94
3	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
4	Heneicosane	296 C21H44	000629-94-7	91
5	2-methyloctacosane	408 C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
Data File : BF118402.D
Acq On : 31 Dec 2019 13:44
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Sample : K6463-04
Misc :
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Instrument :
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ClientSampleId :
GP-7-(4-6)

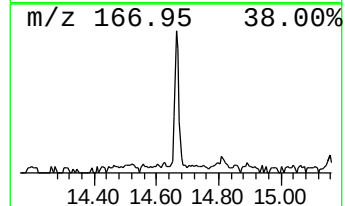
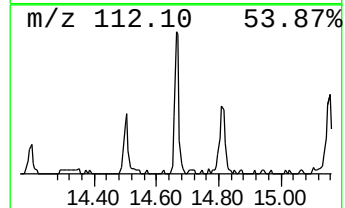
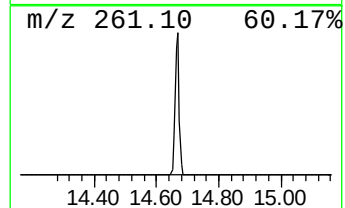
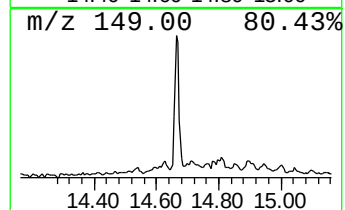
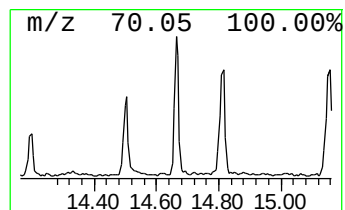
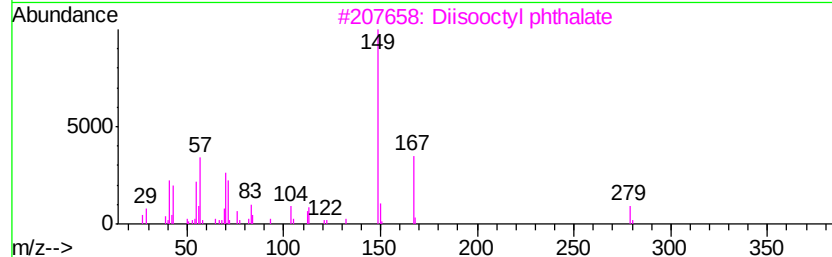
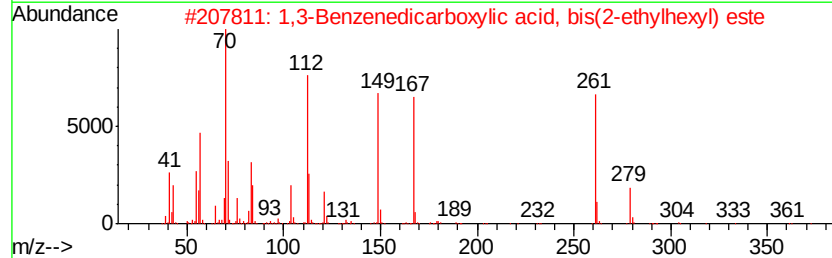
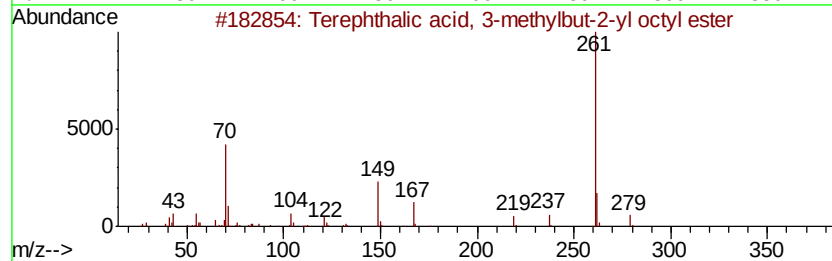
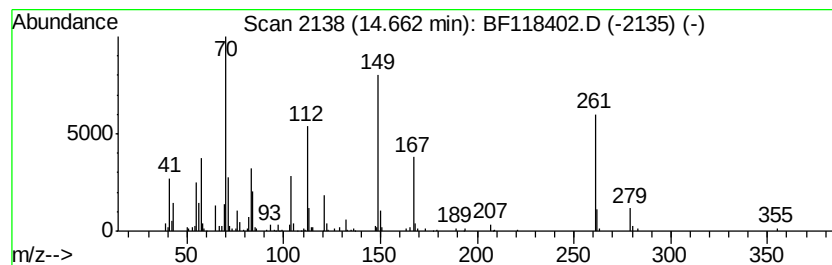
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Terephthalic acid, 3-methyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	3.68 ng	197900	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	64
2		1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	43
3		Diisooctyl phthalate	390	C24H38O4	000131-20-4	38
4		Terephthalic acid, phenyl octyl ...	354	C22H26O4	1000324-06-6	38
5		Terephthalic acid, monochloride,...	296	C16H21ClO3	1000324-22-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
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Acq On : 31 Dec 2019 13:44
Operator : JU
Sample : K6463-04
Misc :
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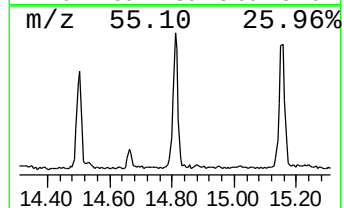
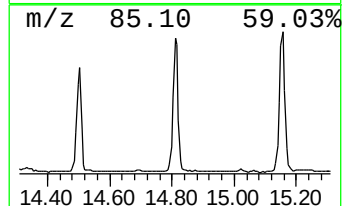
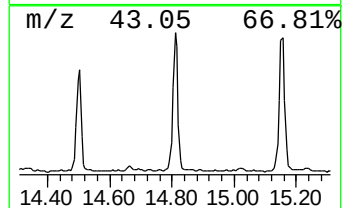
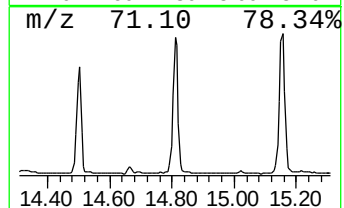
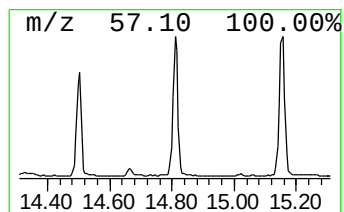
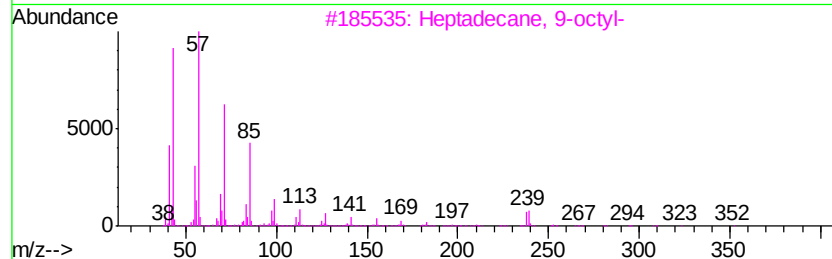
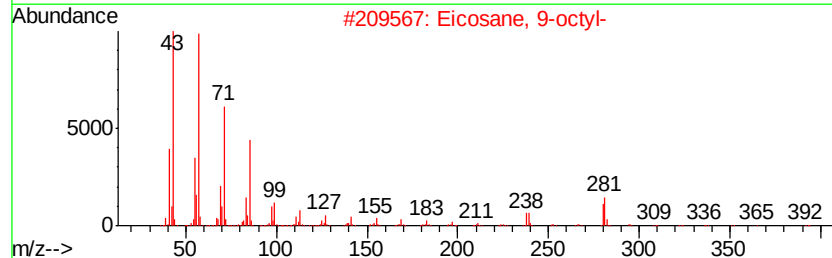
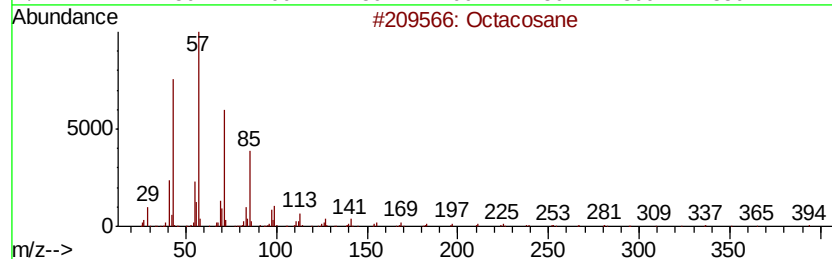
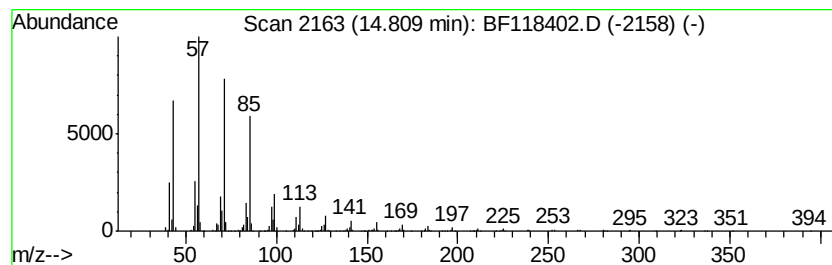
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	10.42 ng	1152760	Perylene-d12	15.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	94
2	Eicosane, 9-octyl-	394 C28H58	013475-77-9	93
3	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
4	Heneicosane	296 C21H44	000629-94-7	91
5	2-methyloctacosane	408 C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
Data File : BF118402.D
Acq On : 31 Dec 2019 13:44
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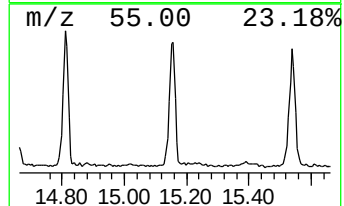
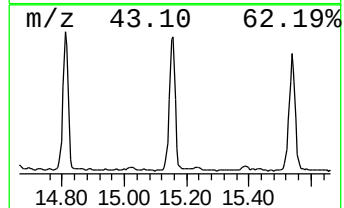
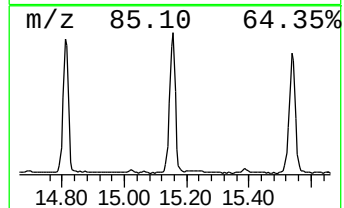
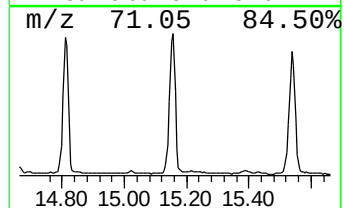
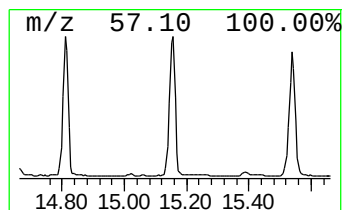
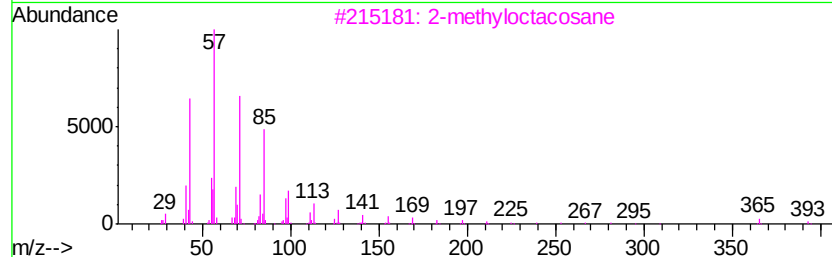
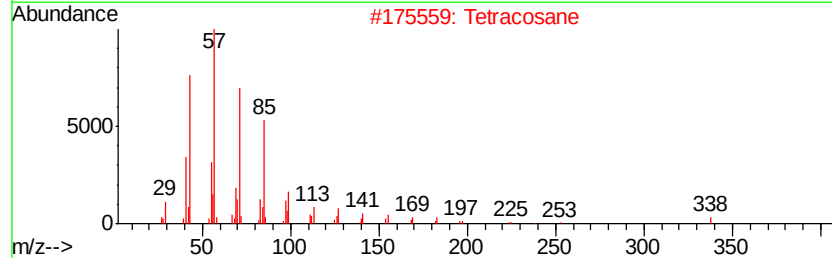
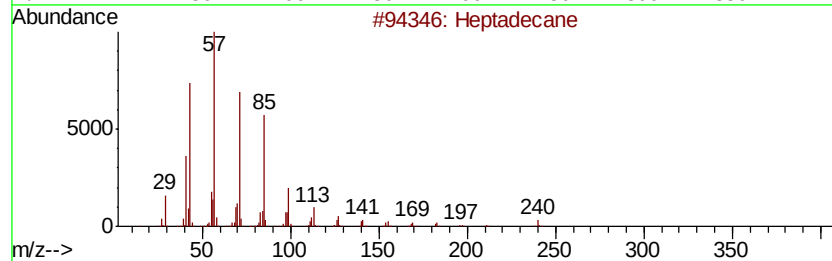
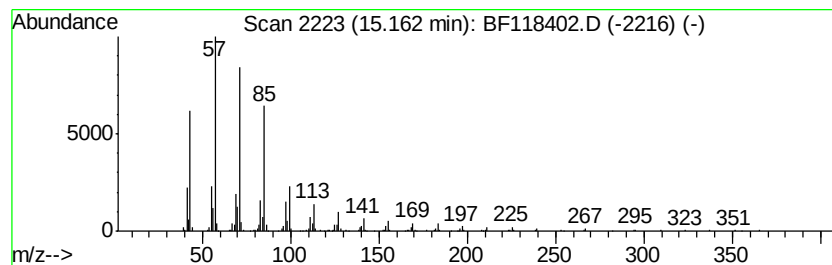
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	11.24 ng	1243520	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	94
2		Tetracosane	338	C24H50	000646-31-1	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
Data File : BF118402.D
Acq On : 31 Dec 2019 13:44
Operator : JU
Sample : K6463-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GP-7-(4-6)

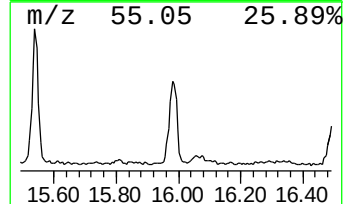
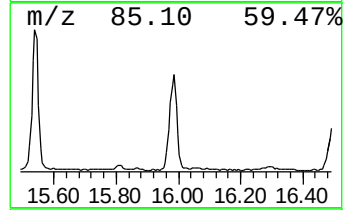
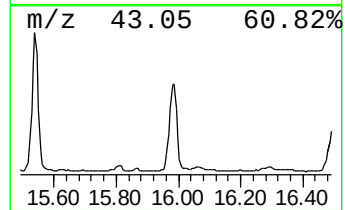
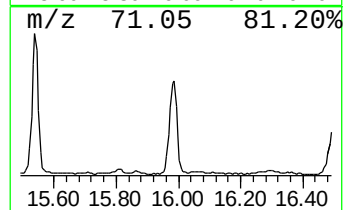
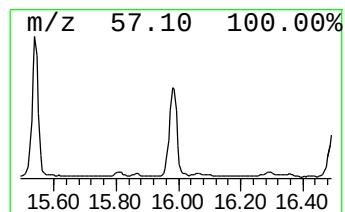
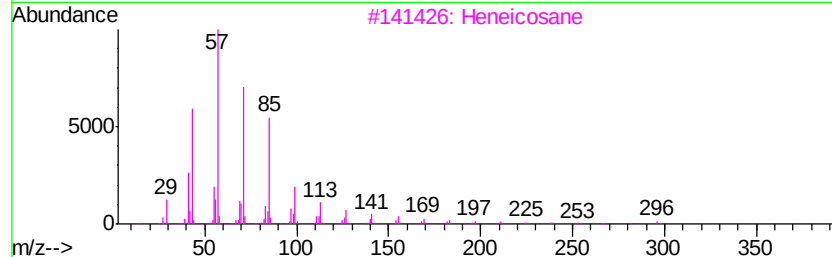
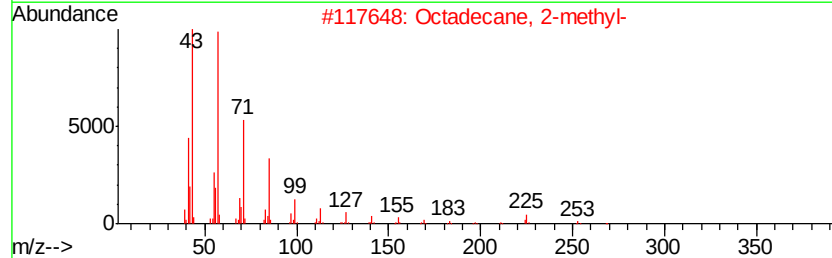
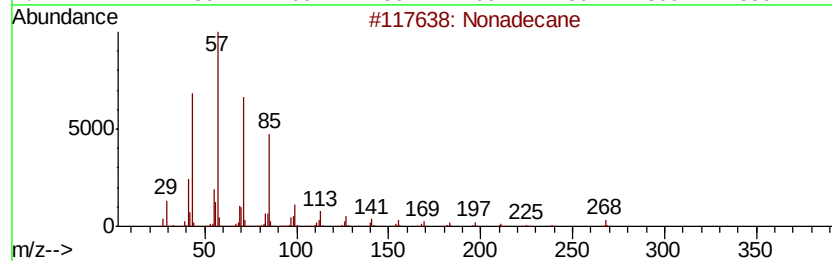
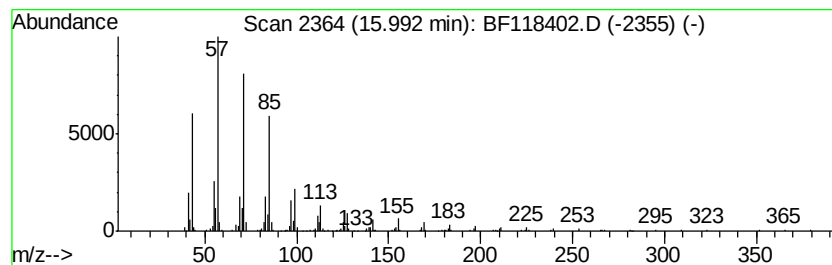
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	9.18 ng	1015360	Perylene-d12	15.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Octadecane, 2-methyl-	268	C19H40	001560-88-9	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123119\
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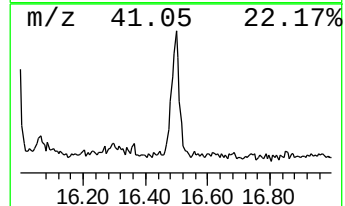
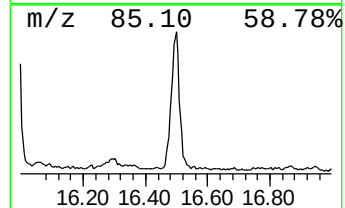
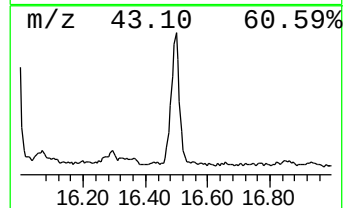
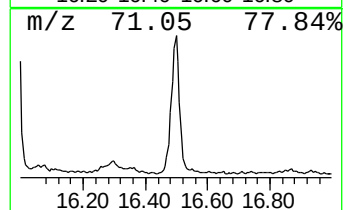
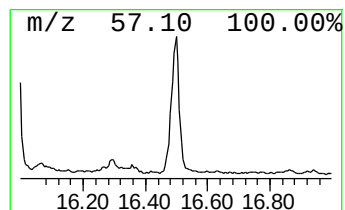
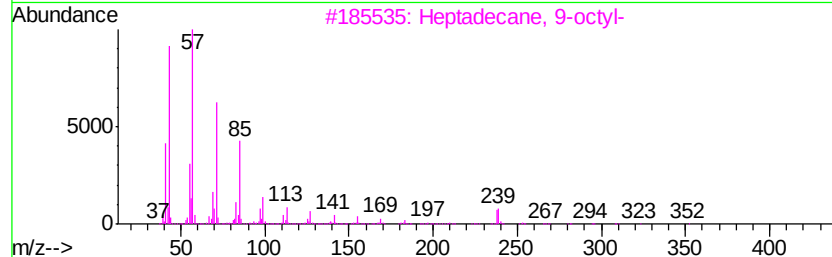
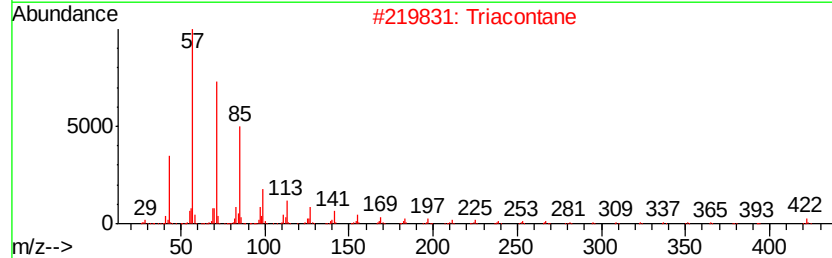
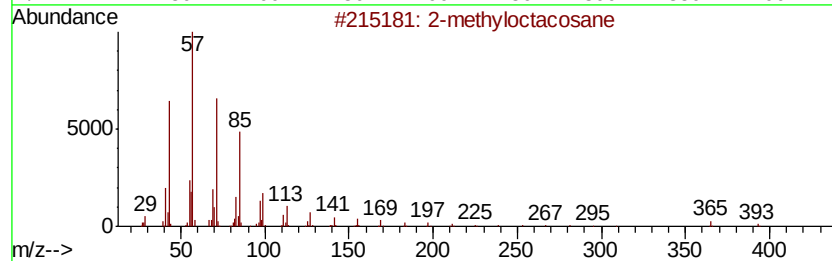
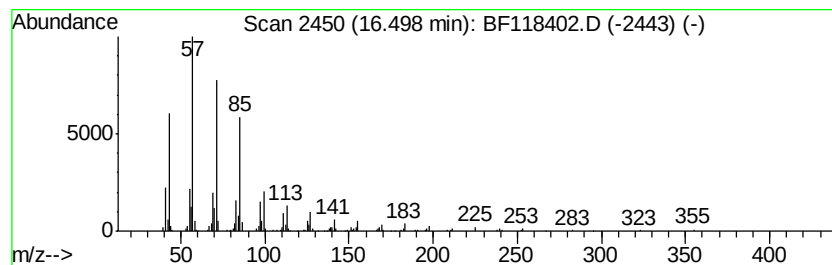
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-methyloctacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	4.75 ng	525538	Perylene-d12	15.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-methyloctacosane	408 C29H60	1000376-72-8	91
2	triacontane	422 C30H62	000638-68-6	91
3	heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
4	pentadecane, 8-hexyl-	296 C21H44	013475-75-7	91
5	heptacosane	380 C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF123119\
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 Sample : K6463-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.06	9.8	ng	354706	1	6.85	719873	20.0
unknown6.62	6.62	68.9	ng	2478420	1	6.85	719873	20.0
n-Hexadecanoic acid	11.92	6.6	ng	408836	4	11.40	1243690	20.0
Octadecanoic acid	12.70	6.0	ng	372953	4	11.40	1243690	20.0
Ethanol, 2-(tetra...	13.88	10.3	ng	555243	5	14.05	1075930	20.0
Heptadecane	14.20	7.9	ng	424495	5	14.05	1075930	20.0
Hexadecane	14.50	14.3	ng	768240	5	14.05	1075930	20.0
Terephthalic acid...	14.66	3.7	ng	197900	5	14.05	1075930	20.0
Octacosane	14.81	10.4	ng	1152760	6	15.55	2213120	20.0
Heneicosane	15.16	11.2	ng	1243520	6	15.55	2213120	20.0
Nonadecane	15.99	9.2	ng	1015360	6	15.55	2213120	20.0
2-methyloctacosane	16.50	4.8	ng	525538	6	15.55	2213120	20.0