

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
 Data File : BF118702.D
 Acq On : 24 Jan 2020 22:59
 Operator : CG/JU
 Sample : L1242-09
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20200062-COMPOSITE-B8

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-BF012020.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	2.140	3	9	17	rVB	1110202	1496787	27.62%	3.367%
2	5.081	501	509	517	rVB2	42337	69222	1.28%	0.156%
3	5.487	573	578	581	rBV	4965140	4692539	86.59%	10.556%
4	6.492	743	749	756	rBV	4846518	4956106	91.46%	11.149%
5	6.863	809	812	815	rBV	1727214	1404920	25.93%	3.160%
6	7.022	831	839	842	rBV	4856561	4132943	76.27%	9.297%
7	7.151	858	861	869	rVB	61645	86129	1.59%	0.194%
8	7.257	875	879	884	rBV	99310	99422	1.83%	0.224%
9	7.422	902	907	910	rBV	3444437	3042213	56.14%	6.843%
10	7.634	939	943	948	rBV	80599	79341	1.46%	0.178%
11	8.145	1026	1030	1033	rBV	2188582	1789432	33.02%	4.025%
12	9.222	1208	1213	1216	rBV	6103154	5418985	100.00%	12.190%
13	9.304	1224	1227	1231	rBV3	63390	85894	1.59%	0.193%
14	9.610	1275	1279	1287	rBV	1026782	892299	16.47%	2.007%
15	9.898	1324	1328	1331	rBV	2264036	1878732	34.67%	4.226%
16	9.928	1331	1333	1337	rVB	85176	73356	1.35%	0.165%
17	10.686	1457	1462	1465	rBV	3177797	2980123	54.99%	6.704%
18	10.816	1480	1484	1491	rVB3	65746	97817	1.81%	0.220%
19	11.380	1576	1580	1583	rBV	1869794	1565656	28.89%	3.522%
20	11.904	1667	1669	1676	rBV2	279080	279410	5.16%	0.629%
21	12.592	1782	1786	1789	rBV	163396	151752	2.80%	0.341%
22	12.686	1799	1802	1810	rBV3	66036	101928	1.88%	0.229%
23	12.821	1822	1825	1827	rVB	99075	88005	1.62%	0.198%
24	12.963	1845	1849	1852	rBV	4238023	3637286	67.12%	8.182%
25	13.539	1945	1947	1954	rVB2	62035	74660	1.38%	0.168%
26	13.857	1998	2001	2007	rBV	380876	447574	8.26%	1.007%
27	14.016	2022	2028	2031	rBV	1603873	1625737	30.00%	3.657%
28	14.180	2053	2056	2061	rVB	134308	137047	2.53%	0.308%
29	14.486	2105	2108	2112	rBV	206826	203133	3.75%	0.457%
30	14.792	2157	2160	2169	rVB2	192784	243247	4.49%	0.547%
31	14.868	2170	2173	2176	rVB	145082	135394	2.50%	0.305%
32	15.051	2201	2204	2211	rVB	81624	180500	3.33%	0.406%
33	15.133	2214	2218	2221	rBV	189853	235049	4.34%	0.529%
34	15.480	2272	2277	2280	rBV	1342392	1659004	30.61%	3.732%

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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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012020.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.510	2280	2282	2288	rVB	197599	239125	4.41%	0.538%
36	15.951	2353	2357	2361	rVB	122223	174167	3.21%	0.392%

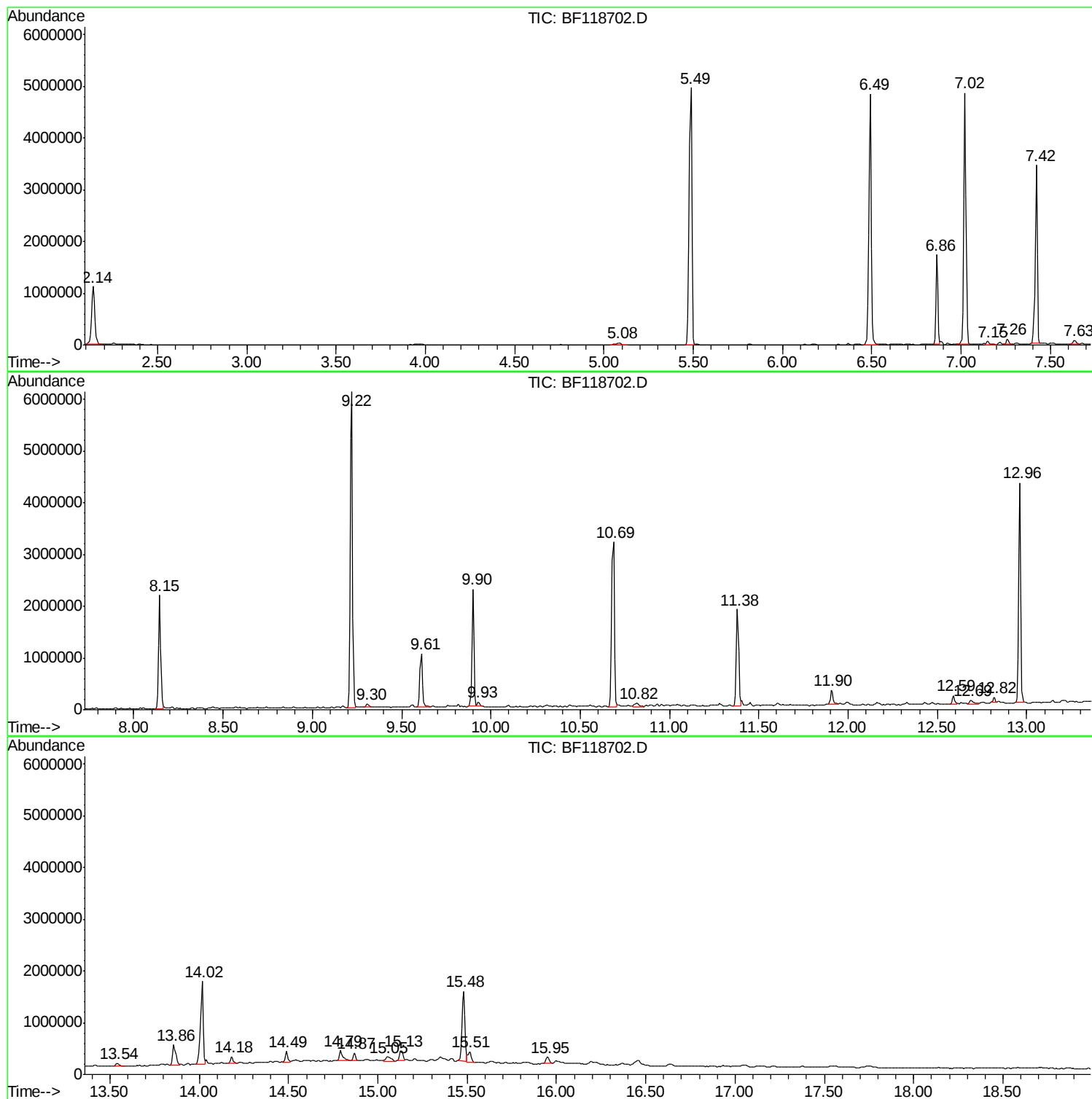
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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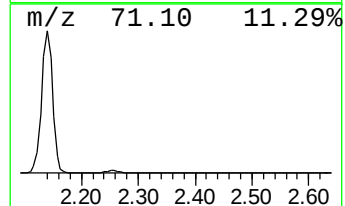
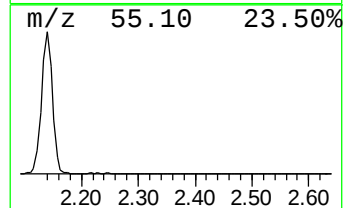
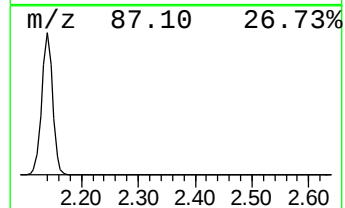
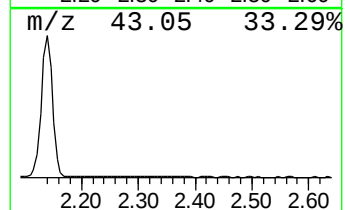
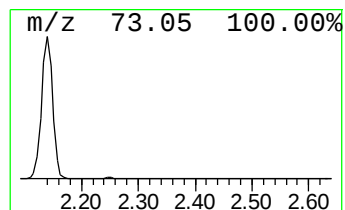
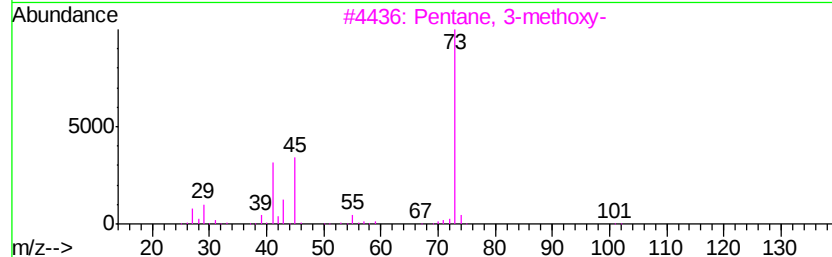
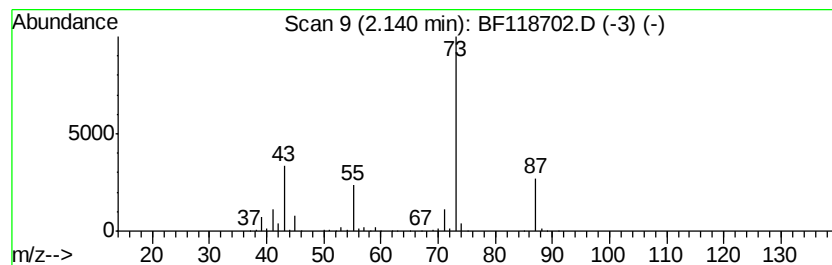
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.14	21.31 ng	1496790	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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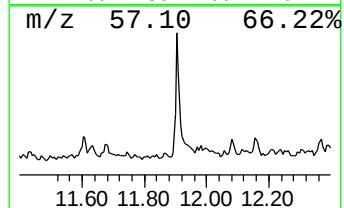
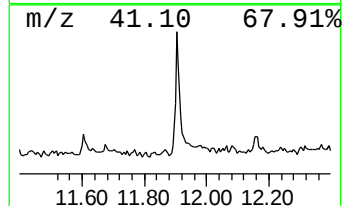
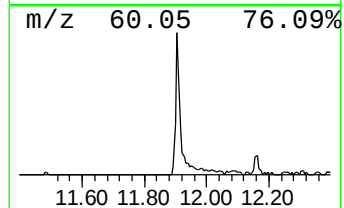
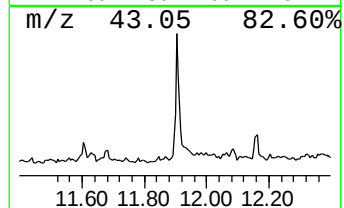
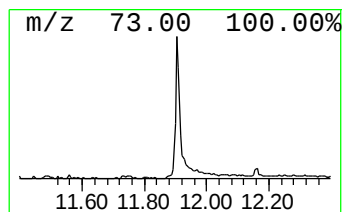
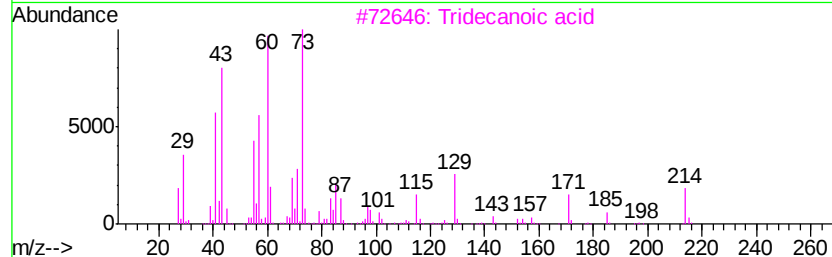
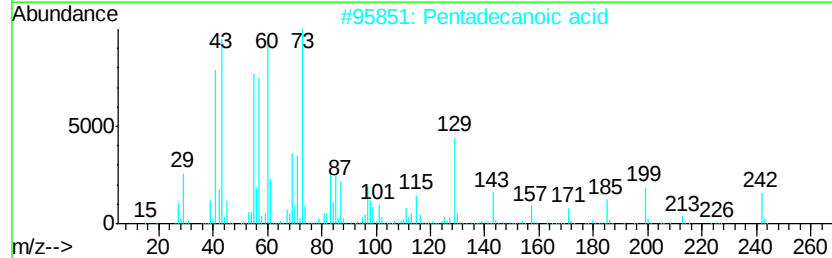
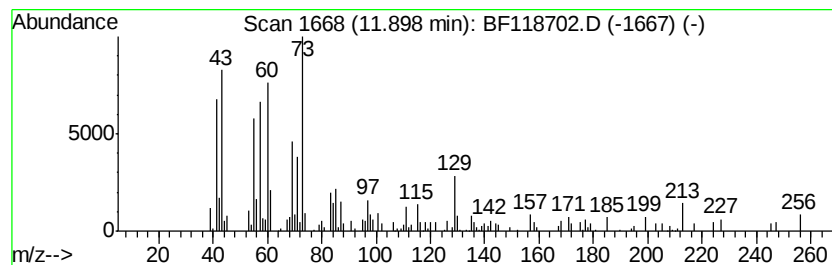
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.90	3.57 ng	279410	Phenanthrene-d10	11.38	
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	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	70
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	68
5	Dodecanoic acid	200	C12H24O2	000143-07-7	58



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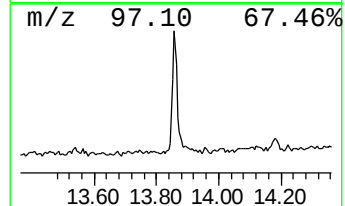
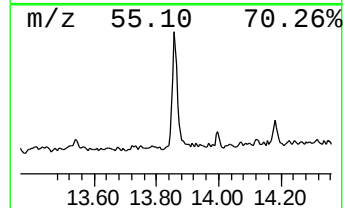
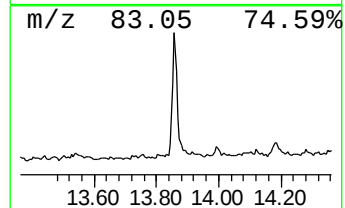
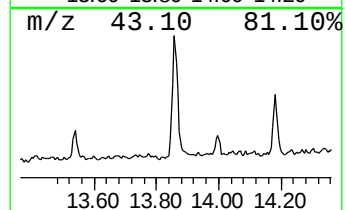
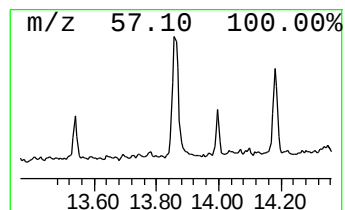
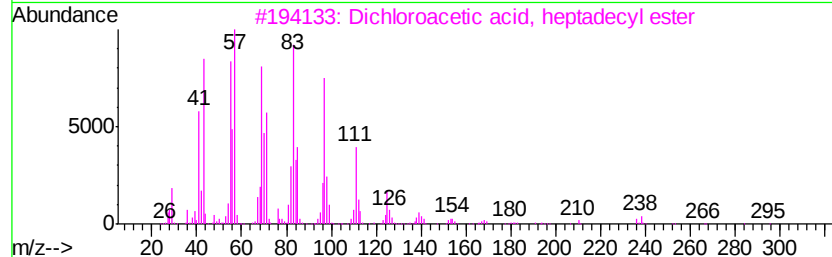
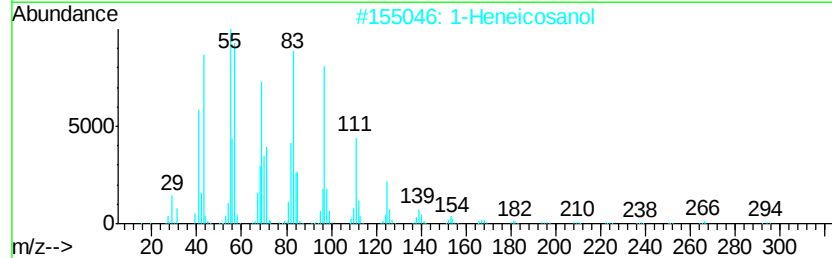
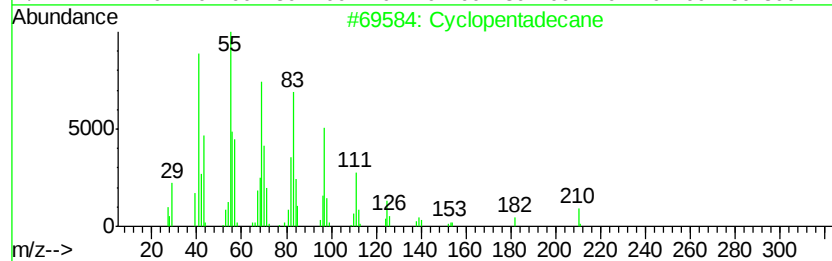
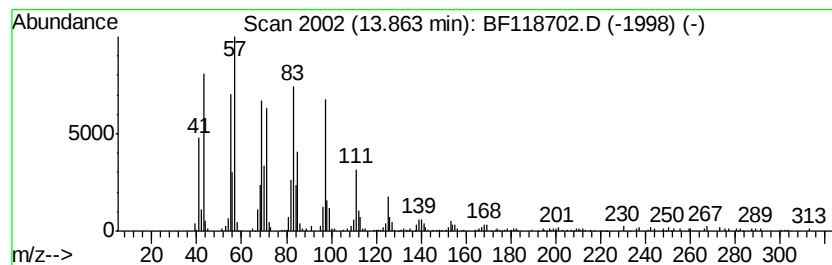
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopentadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	5.51 ng	447574	Chrysene-d12	14.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentadecane	210	C15H30	000295-48-7	97
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
5		2-Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91



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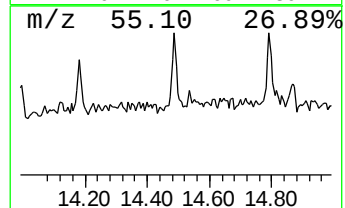
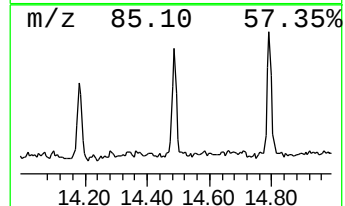
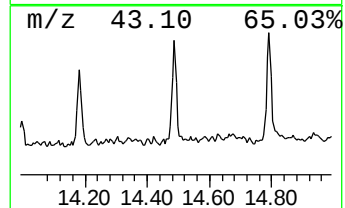
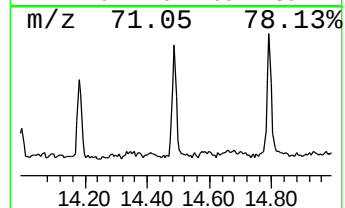
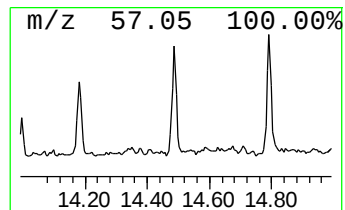
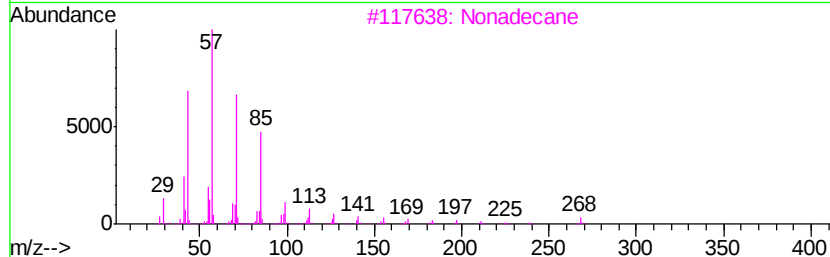
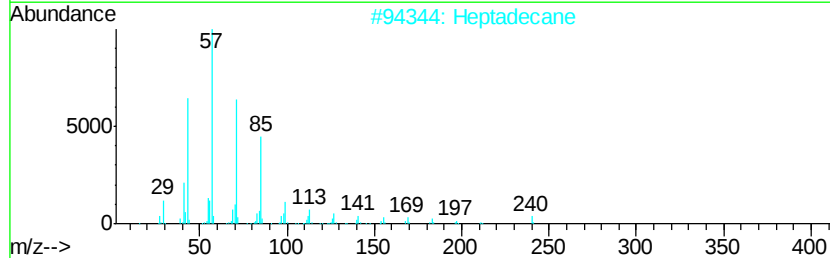
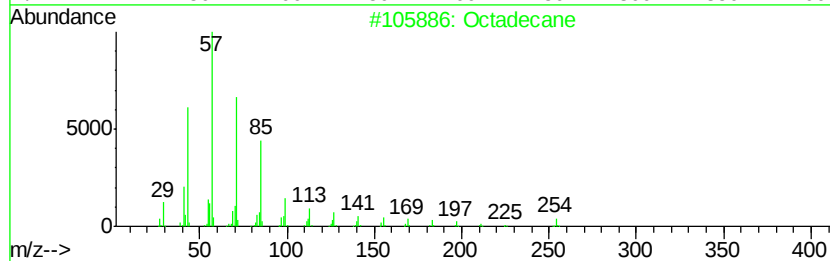
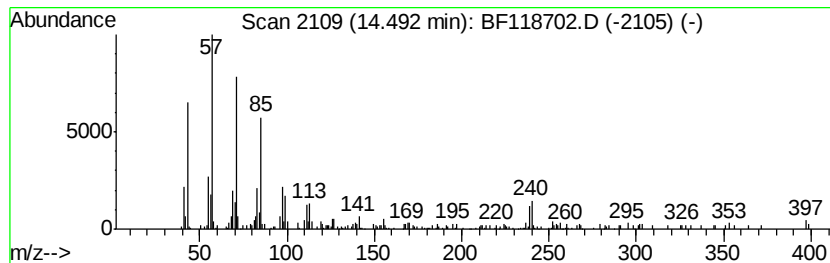
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.49	2.50 ng	203133	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	95
2	Heptadecane	240	C17H36	000629-78-7	95
3	Nonadecane	268	C19H40	000629-92-5	93
4	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
5	Tetratetracontane	619	C44H90	007098-22-8	91



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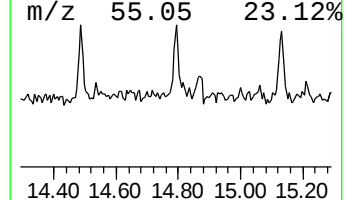
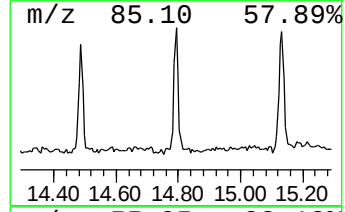
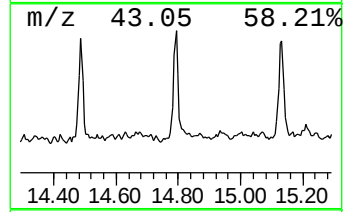
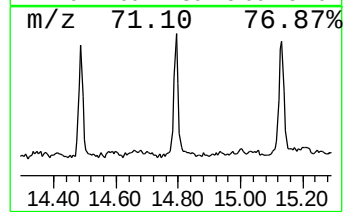
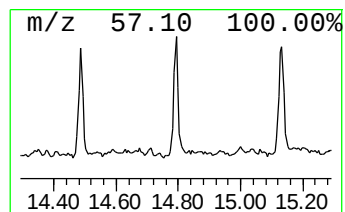
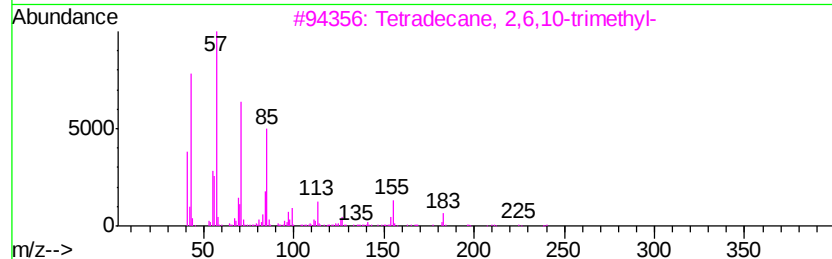
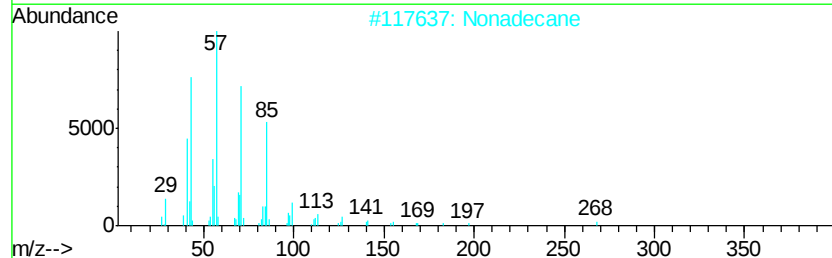
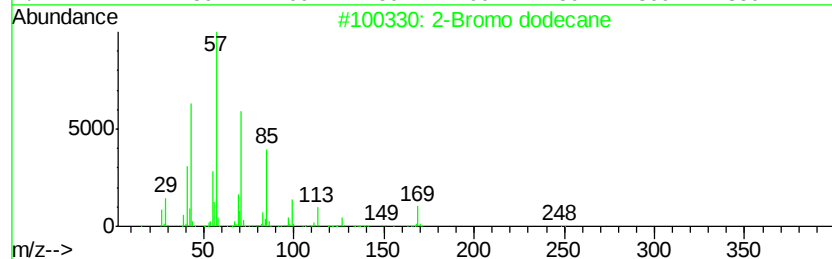
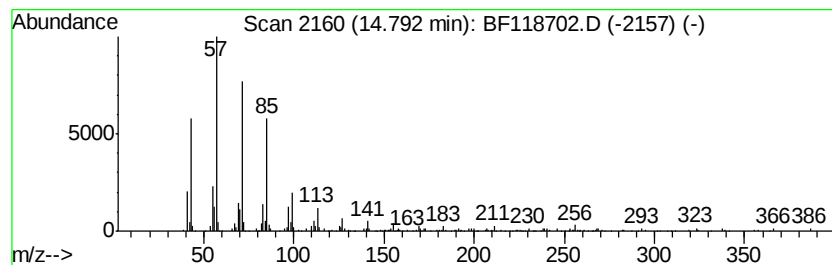
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Bromo dodecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	2.93 ng	243247	Perylene-d12	15.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
2		Nonadecane	268	C19H40	000629-92-5	91
3		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



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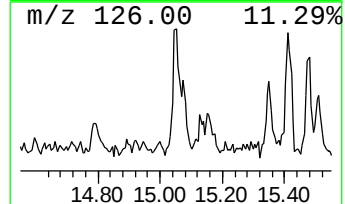
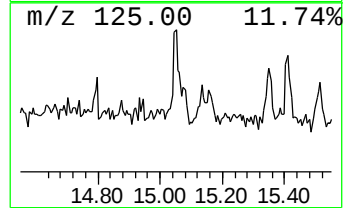
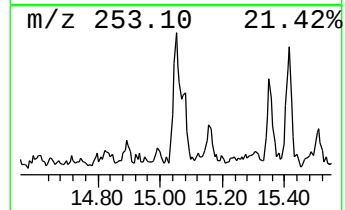
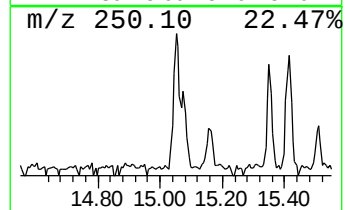
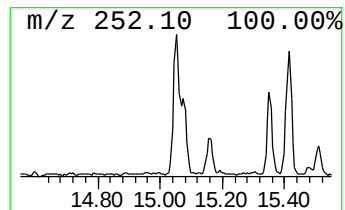
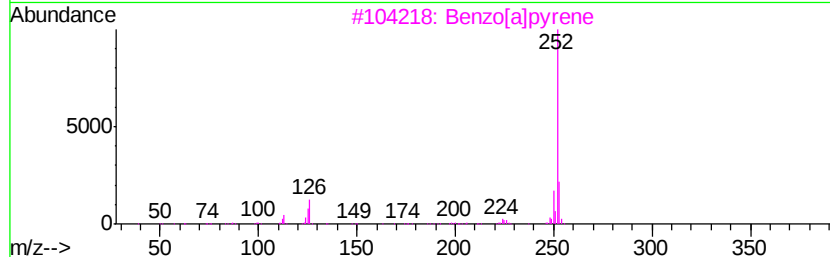
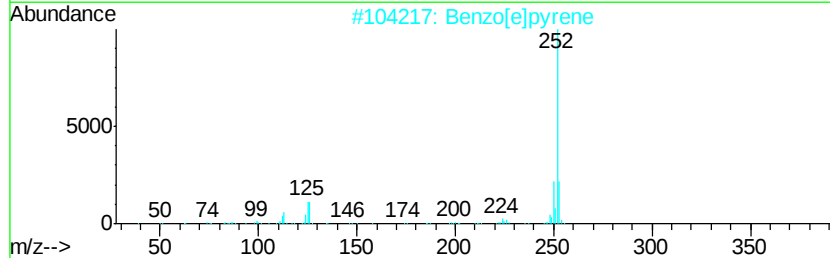
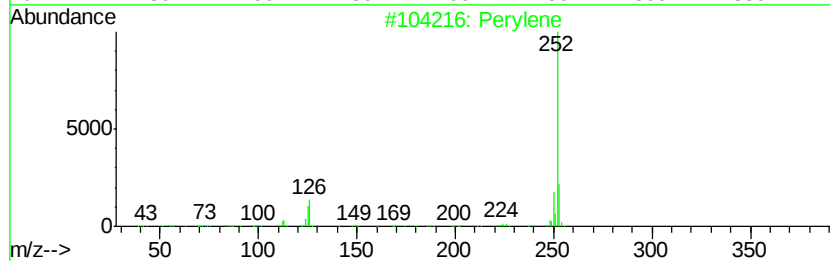
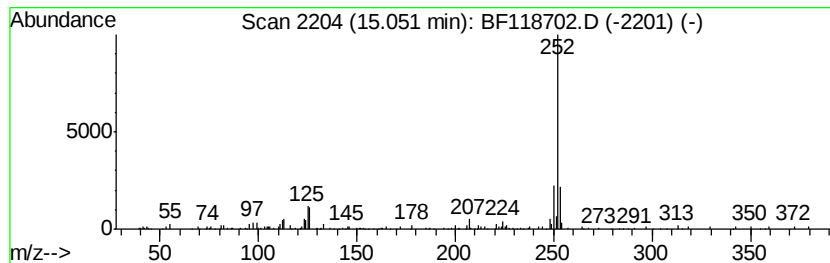
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Perylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	2.18 ng	180500	Perylene-d12	15.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene		252	C20H12	000198-55-0	95
2	Benzo[e]pyrene		252	C20H12	000192-97-2	93
3	Benzo[a]pyrene		252	C20H12	000050-32-8	93
4	Benzo[e]acephenanthrylene		252	C20H12	000205-99-2	91
5	Benzo[j]fluoranthene		252	C20H12	000205-82-3	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
Data File : BF118702.D
Acq On : 24 Jan 2020 22:59
Operator : CG/JU
Sample : L1242-09
Misc :
ALS Vial : 22 Sample Multiplier: 1

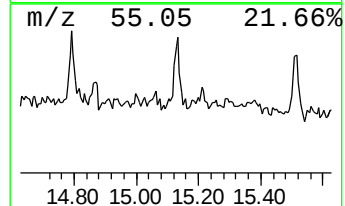
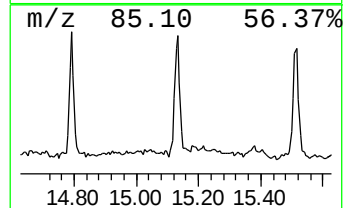
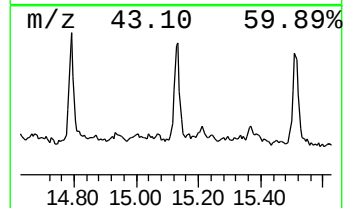
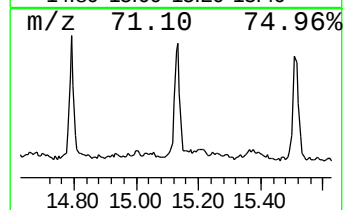
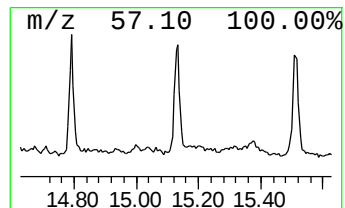
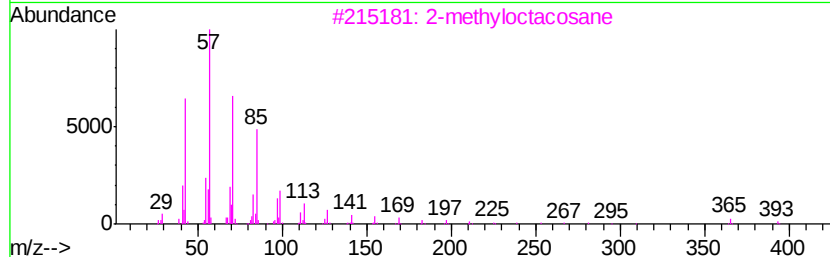
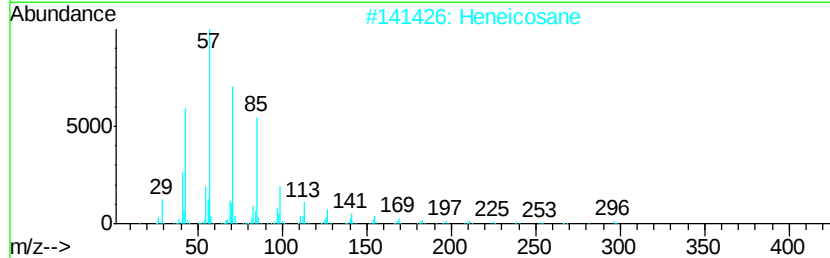
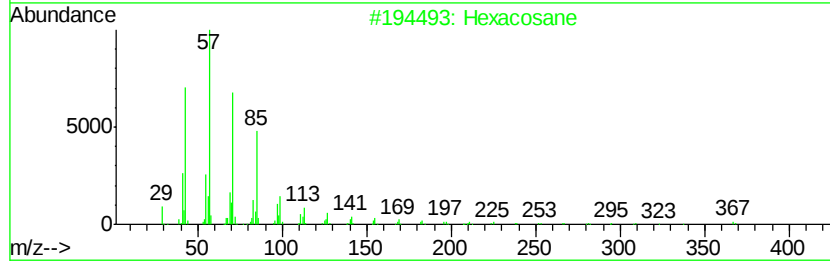
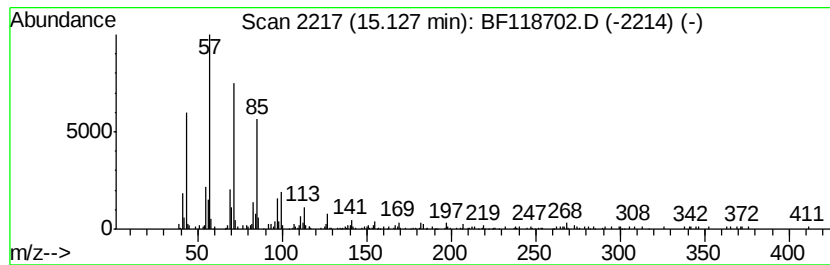
Instrument :
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ClientSampleId :
20200062-COMPOSITE-B8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.13	2.83 ng	235049	Perylene-d12	15.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	91
2	Heneicosane	296 C21H44	000629-94-7	91
3	2-methyloctacosane	408 C29H60	1000376-72-8	91
4	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
5	Triacontane	422 C30H62	000638-68-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
Data File : BF118702.D
Acq On : 24 Jan 2020 22:59
Operator : CG/JU
Sample : L1242-09
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20200062-COMPOSITE-B8

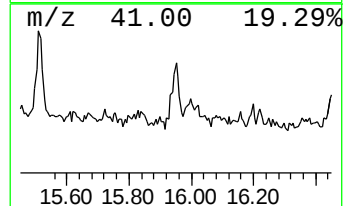
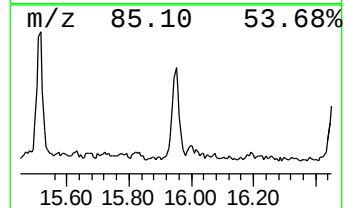
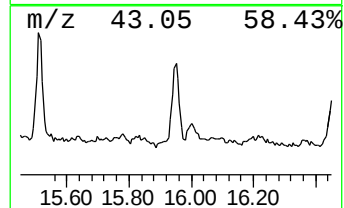
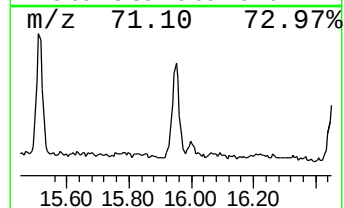
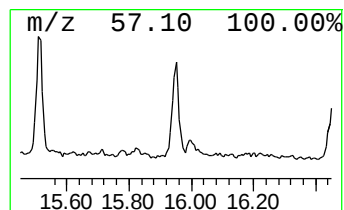
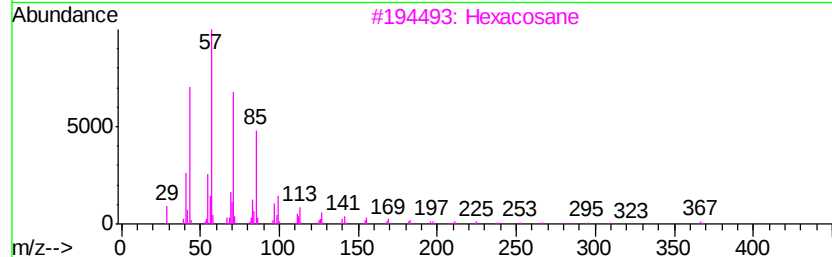
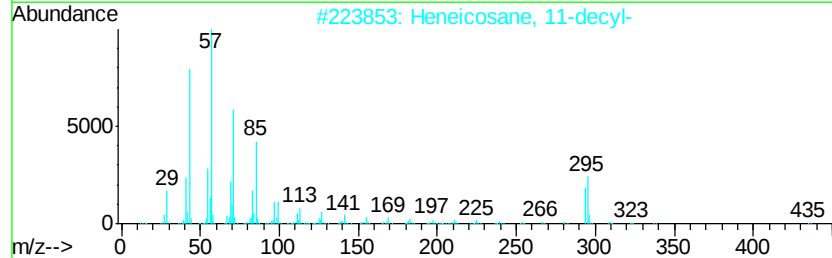
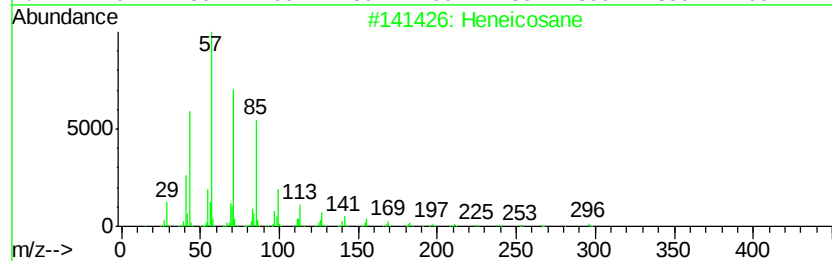
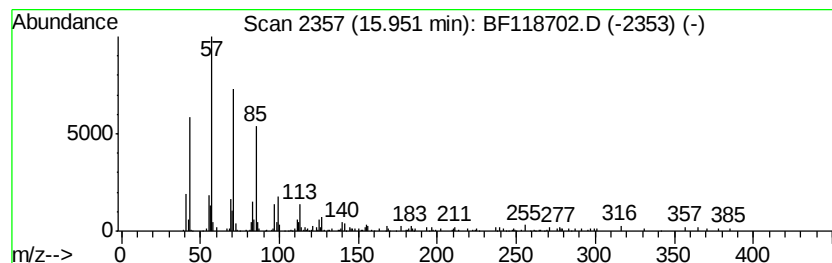
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	2.10 ng	174167	Perylene-d12	15.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	90
2	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	90
3	Hexacosane		366	C26H54	000630-01-3	90
4	Dodecane, 2-methyl-		184	C13H28	001560-97-0	90
5	Tetratetracontane		619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012420\
Data File : BF118702.D
Acq On : 24 Jan 2020 22:59
Operator : CG/JU
Sample : L1242-09
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
20200062-COMPOSITE-B8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012020.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...	2.14	21.3	ng	1496790	1	6.86	1404920	20.0
n-Hexadecanoic acid	11.90	3.6	ng	279410	4	11.38	1565660	20.0
Cyclopentadecane	13.86	5.5	ng	447574	5	14.02	1625740	20.0
Octadecane	14.49	2.5	ng	203133	5	14.02	1625740	20.0
2-Bromo dodecane	14.79	2.9	ng	243247	6	15.48	1659000	20.0
Perylene	15.05	2.2	ng	180500	6	15.48	1659000	20.0
Hexacosane	15.13	2.8	ng	235049	6	15.48	1659000	20.0
Heneicosane	15.95	2.1	ng	174167	6	15.48	1659000	20.0