

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
 Data File : BF117858.D
 Acq On : 19 Nov 2019 14:42
 Operator : JU
 Sample : K5864-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01-COMP

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-BF111319.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	2.187	11	17	24	rVB	566424	738606	14.55%	1.600%
2	5.128	507	517	525	rBV	573663	689484	13.58%	1.494%
3	5.528	580	585	588	rBV	3870087	3501379	68.97%	7.585%
4	6.534	751	756	770	rBV	3842335	3986765	78.54%	8.637%
5	6.681	777	781	784	rBV	4447089	3880708	76.45%	8.407%
6	6.916	817	821	824	rBV	1583524	1259858	24.82%	2.729%
7	7.075	844	848	851	rBV	3325659	3042932	59.94%	6.592%
8	7.475	911	916	919	rBV	2771555	2384641	46.98%	5.166%
9	8.204	1036	1040	1043	rBV	1882861	1667017	32.84%	3.611%
10	9.281	1218	1223	1226	rBV	5684845	4751013	93.59%	10.292%
11	9.675	1286	1290	1293	rBV	490652	418675	8.25%	0.907%
12	9.963	1335	1339	1343	rBV	2360360	2031252	40.01%	4.400%
13	10.757	1469	1474	1477	rBV	3117931	2861458	56.37%	6.199%
14	11.457	1589	1593	1596	rBV	2611125	2156252	42.48%	4.671%
15	11.522	1599	1604	1608	rVB4	52841	62307	1.23%	0.135%
16	11.669	1626	1629	1632	rBV	51848	55691	1.10%	0.121%
17	11.975	1673	1681	1686	rBV2	219558	272087	5.36%	0.589%
18	12.674	1797	1800	1809	rBV	105547	146334	2.88%	0.317%
19	12.769	1811	1816	1822	rBV	346691	431959	8.51%	0.936%
20	12.904	1834	1839	1840	rBV2	73695	72665	1.43%	0.157%
21	12.927	1840	1843	1856	rVB	257830	344489	6.79%	0.746%
22	13.045	1859	1863	1867	rBV	5221415	5076385	100.00%	10.997%
23	13.957	2014	2018	2036	rVB4	257682	604780	11.91%	1.310%
24	14.104	2037	2043	2046	rBV	2003578	1751104	34.50%	3.793%
25	14.268	2067	2071	2073	rBV	77609	79637	1.57%	0.173%
26	14.574	2119	2123	2125	rBV	145358	157741	3.11%	0.342%
27	14.886	2173	2176	2181	rVB	172279	182794	3.60%	0.396%
28	15.245	2233	2237	2241	rVB	196212	212651	4.19%	0.461%
29	15.610	2294	2299	2303	rBV	1197629	1517058	29.88%	3.286%
30	15.645	2303	2305	2310	rVB	186428	207053	4.08%	0.449%
31	16.104	2379	2383	2389	rVB	141272	214047	4.22%	0.464%
32	16.639	2470	2474	2479	rVB2	61237	98751	1.95%	0.214%
33	17.745	2658	2662	2669	rVB9	66930	128875	2.54%	0.279%
34	18.033	2703	2711	2732	rBV9	244104	1175086	23.15%	2.546%

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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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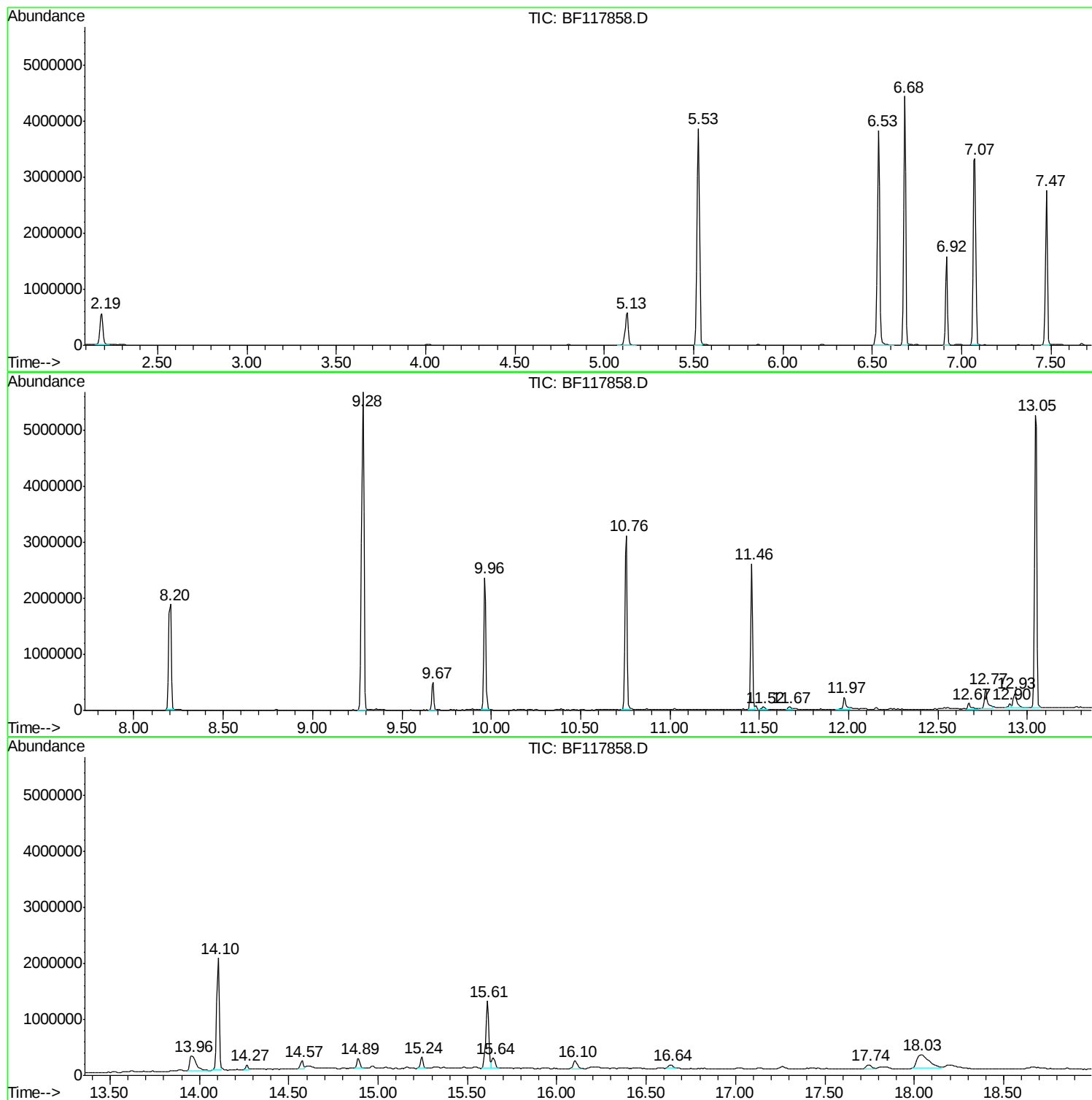
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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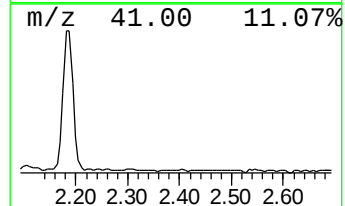
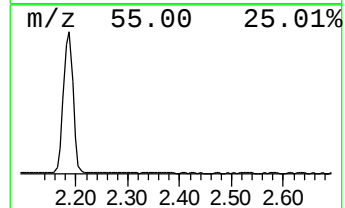
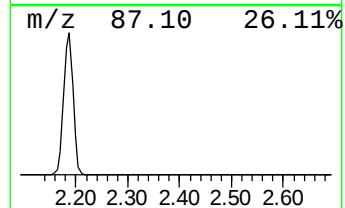
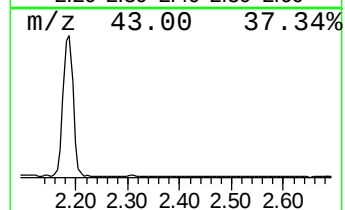
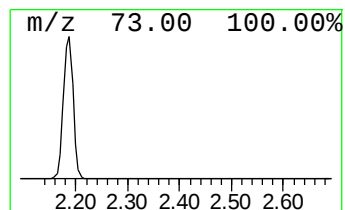
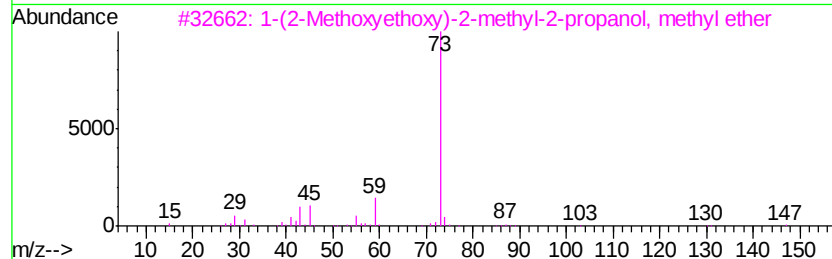
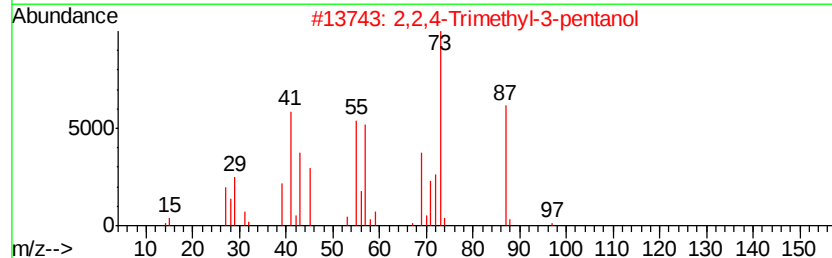
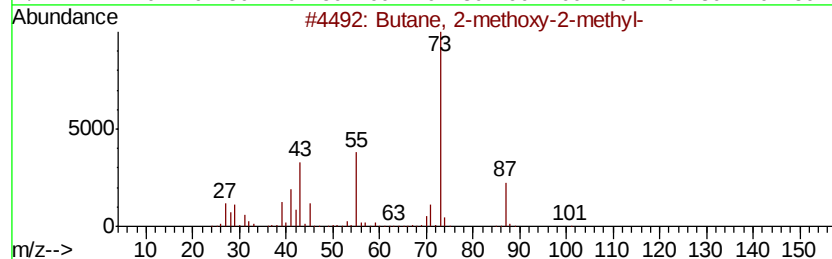
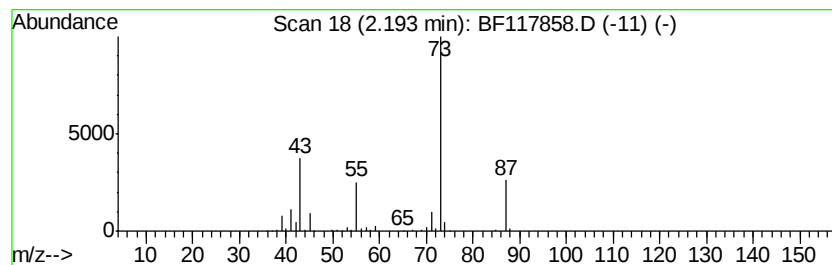
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	11.73 ng	738606	1,4-Dichlorobenzene-d4	6.92

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		1-(2-Methoxyethoxy)-2-methyl-2-propanol	162	C8H18O3	1000364-24-4	25
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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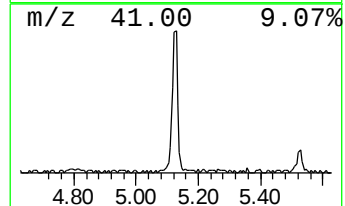
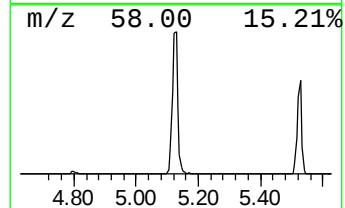
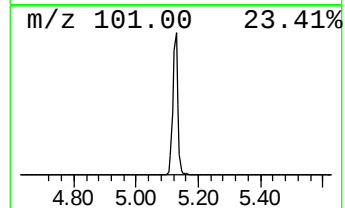
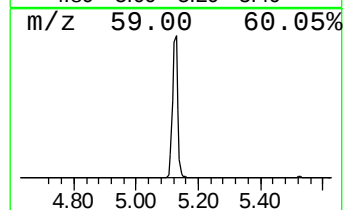
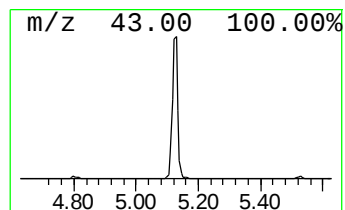
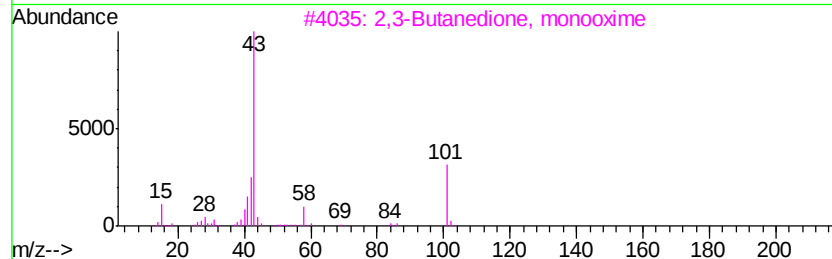
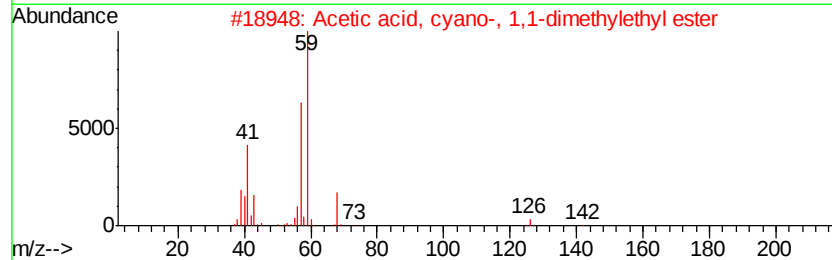
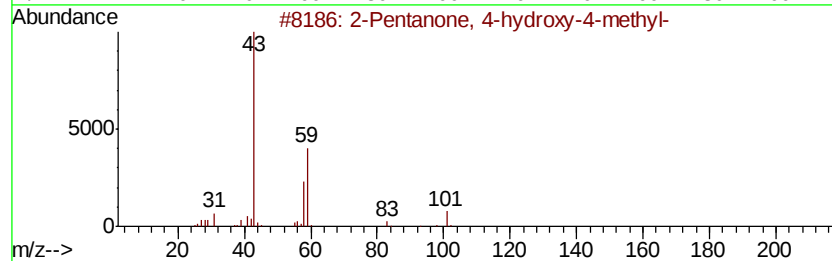
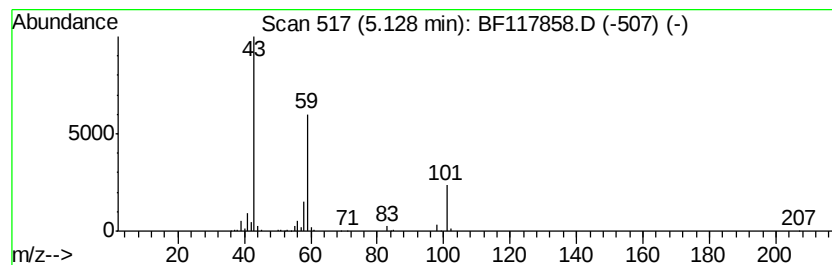
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	10.95 ng	689484	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9



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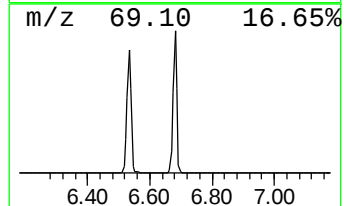
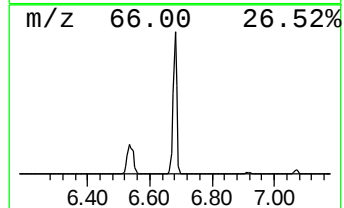
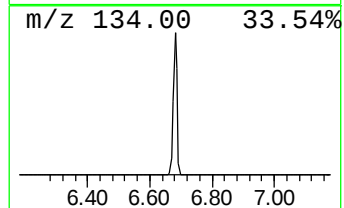
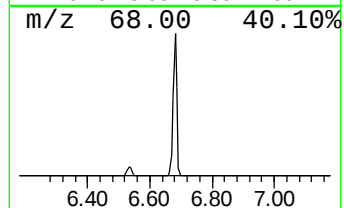
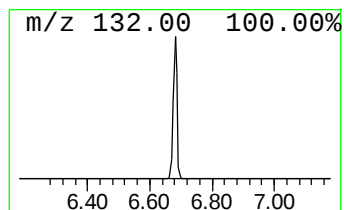
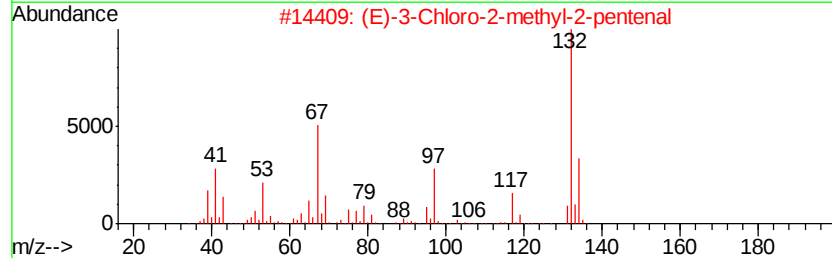
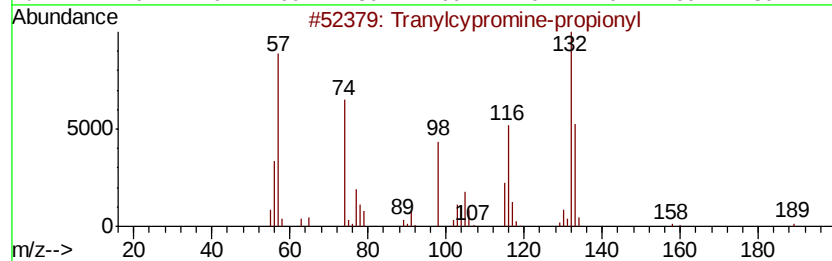
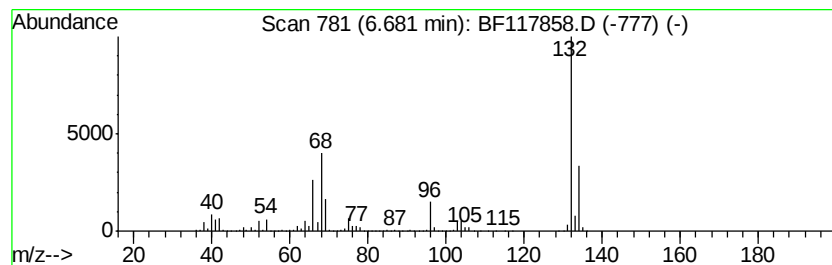
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	61.61 ng	3880710	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Bicvclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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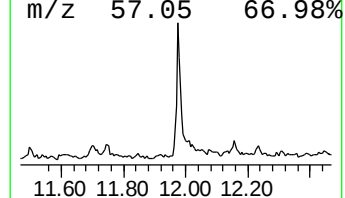
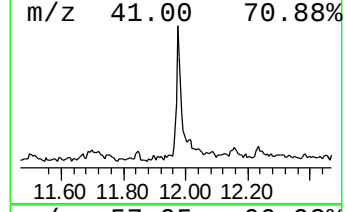
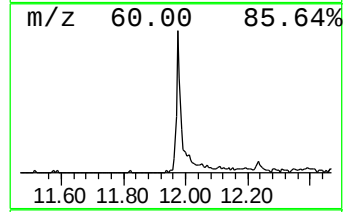
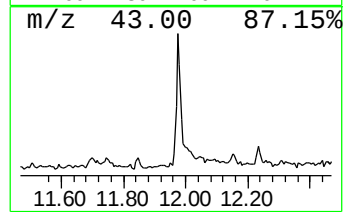
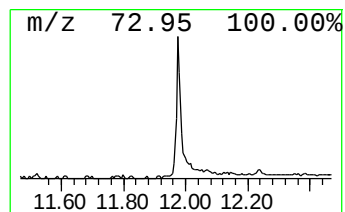
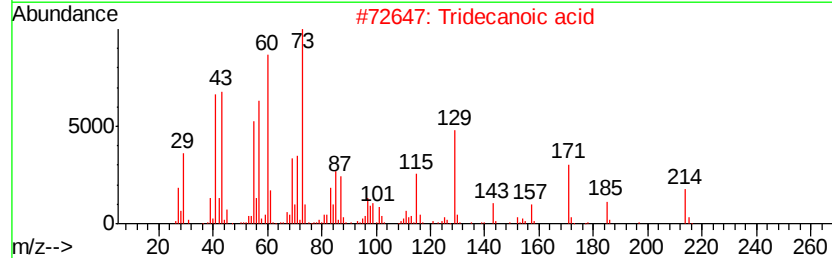
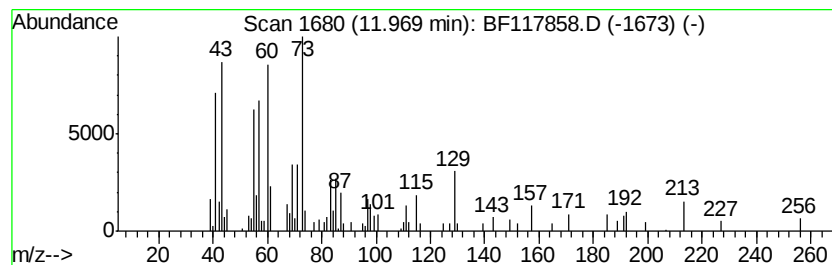
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.52 ng	272087	Phenanthrene-d10	11.46

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		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	41



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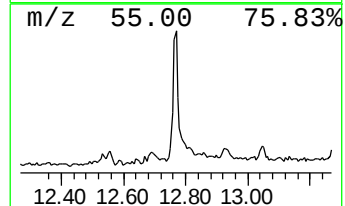
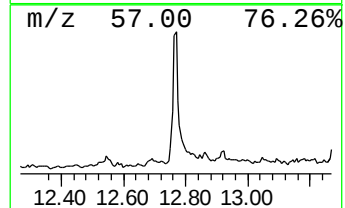
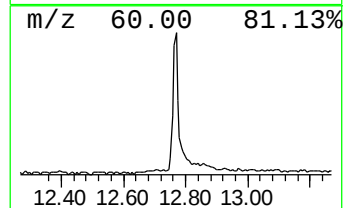
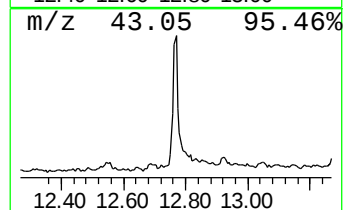
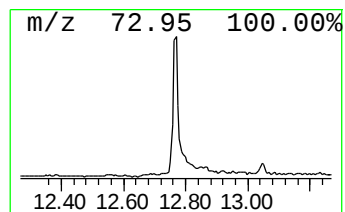
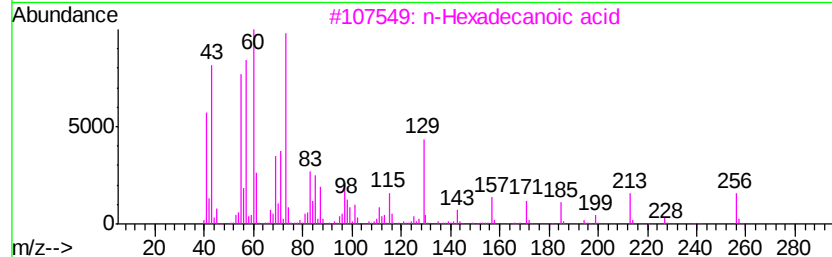
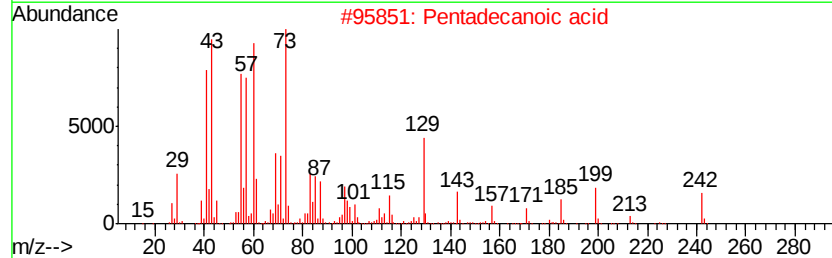
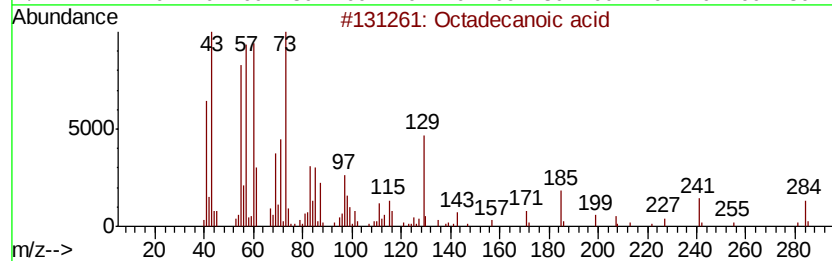
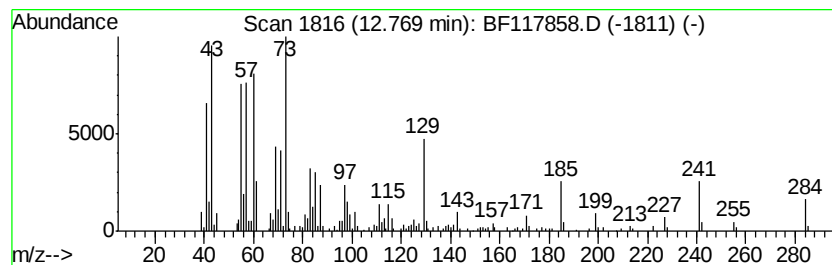
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.77	4.01 ng	431959	Phenanthrene-d10		11.46
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	Tridecanoic acid	214	C13H26O2	000638-53-9	74



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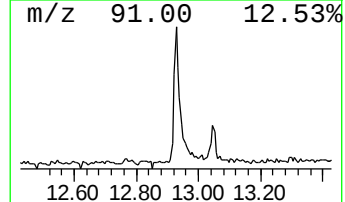
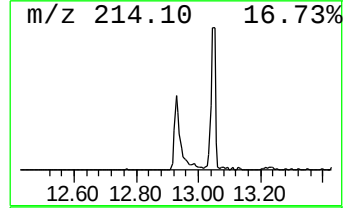
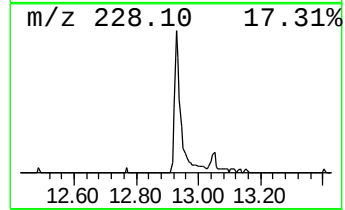
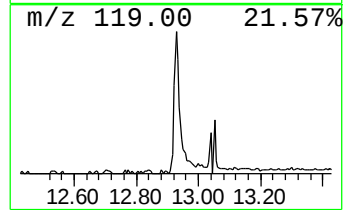
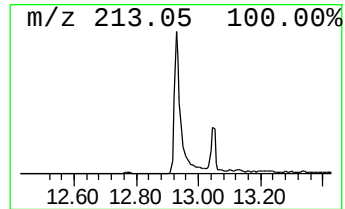
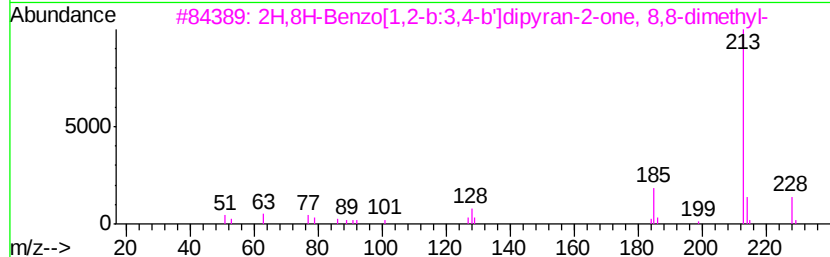
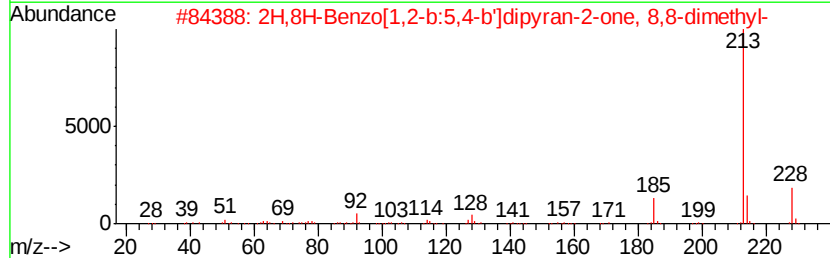
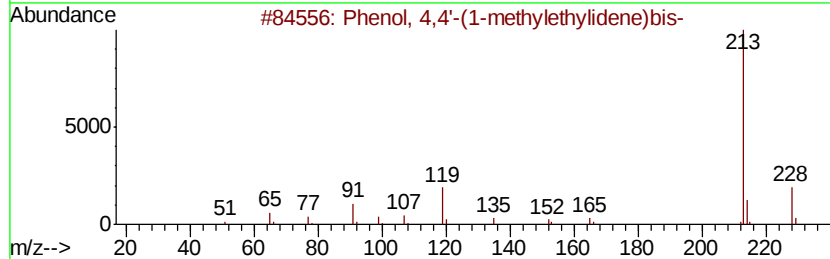
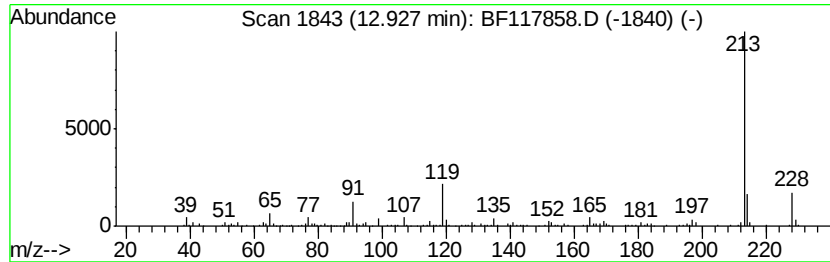
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenol, 4,4'-(1-methylethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.93	3.93 ng	344489	Chrysene-d12	14.10		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98	
2	2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	53	
3	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	53	
4	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	52	
5	As-Indacen-1(2H)-one, 3,6,7,8-te...	228	C16H20O	055591-18-9	42	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117858.D
Acq On : 19 Nov 2019 14:42
Operator : JU
Sample : K5864-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-01-COMP

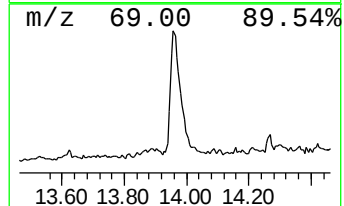
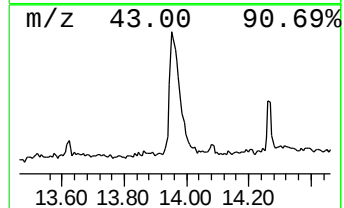
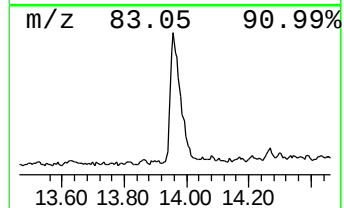
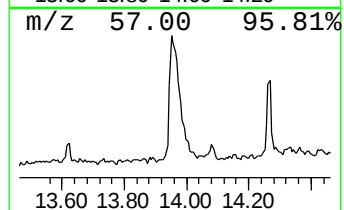
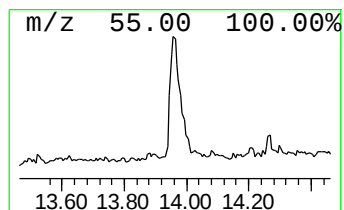
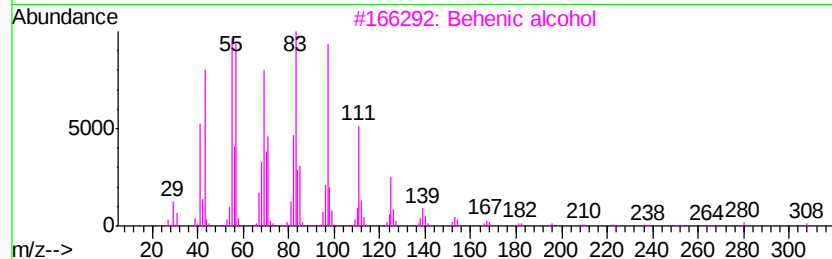
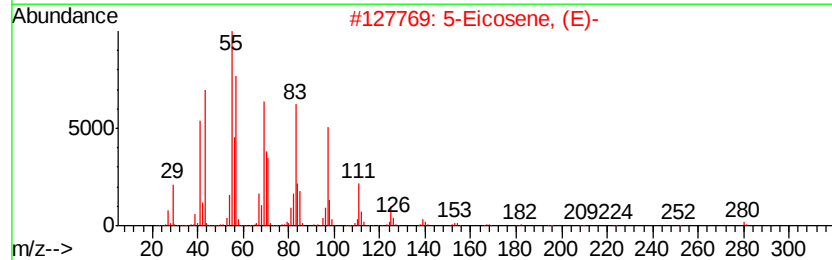
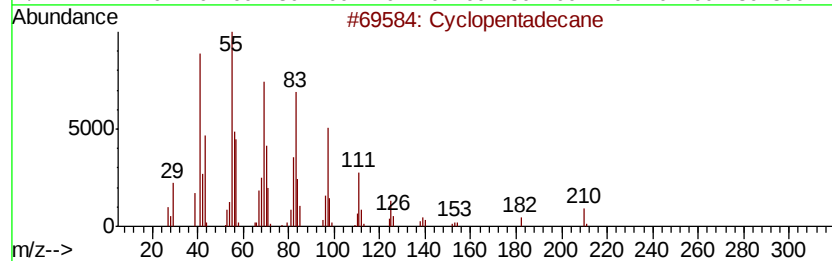
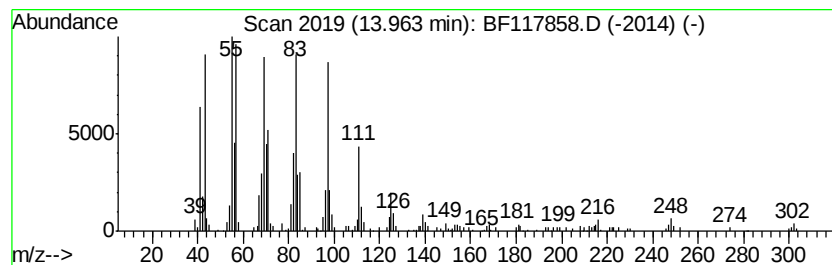
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclopentadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	6.91 ng	604780	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	98
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	96
3		Behenic alcohol	326	C22H46O	000661-19-8	95
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	95
5		1-Heneicosanol	312	C21H44O	015594-90-8	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117858.D
Acq On : 19 Nov 2019 14:42
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Sample : K5864-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

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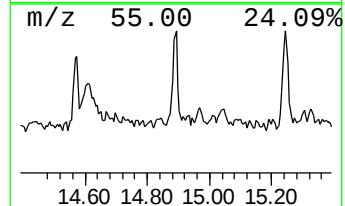
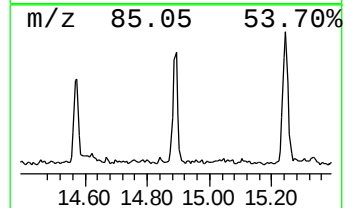
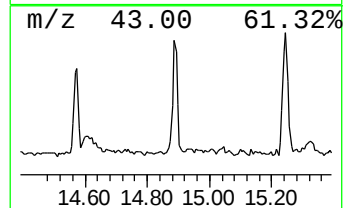
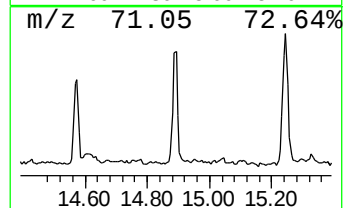
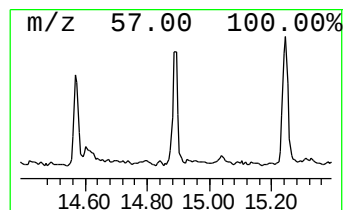
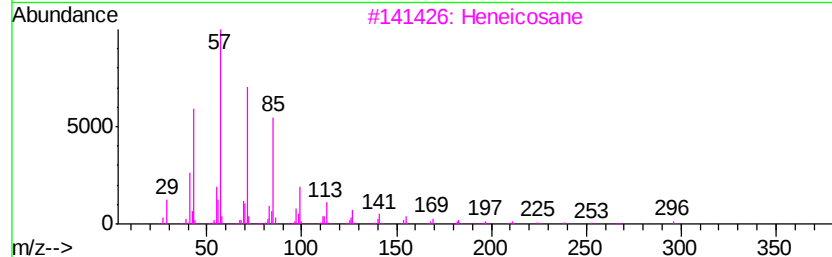
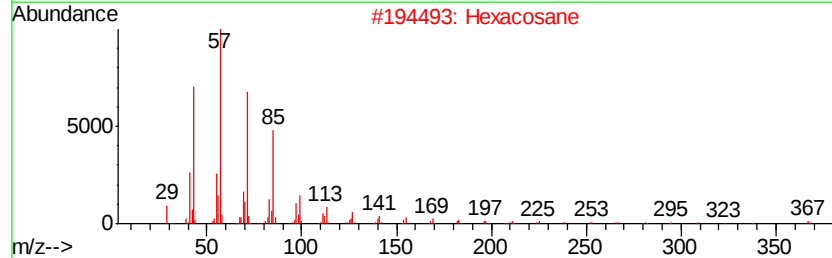
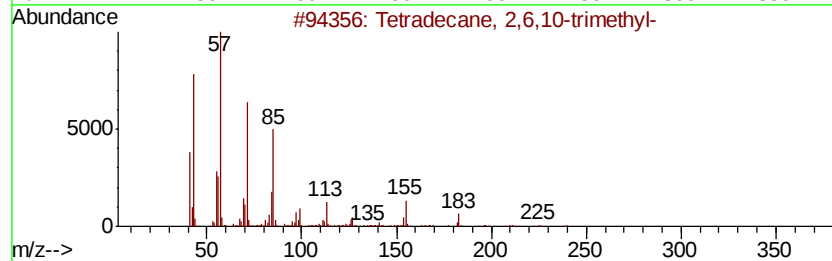
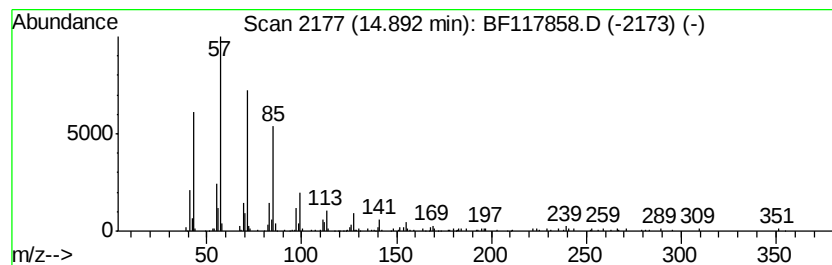
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetradecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	2.41 ng	182794	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87
5		Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117858.D
Acq On : 19 Nov 2019 14:42
Operator : JU
Sample : K5864-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

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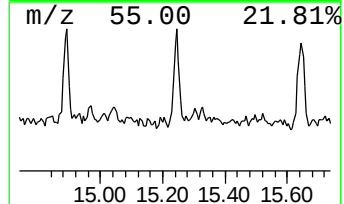
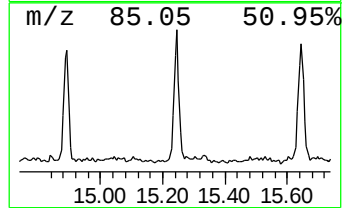
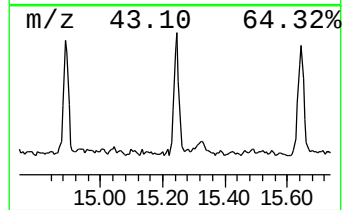
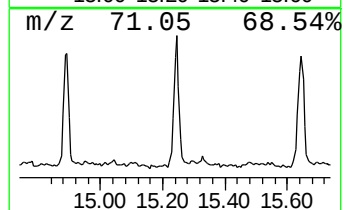
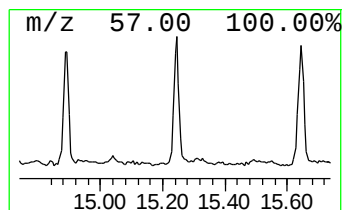
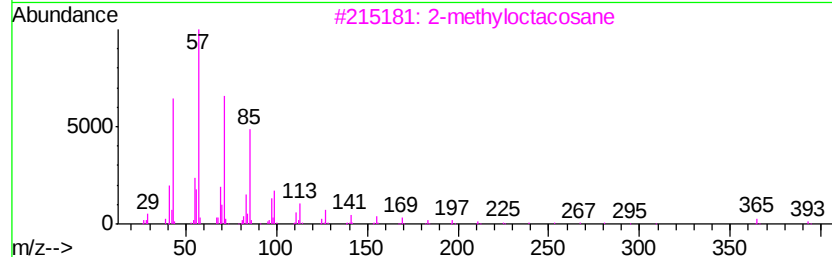
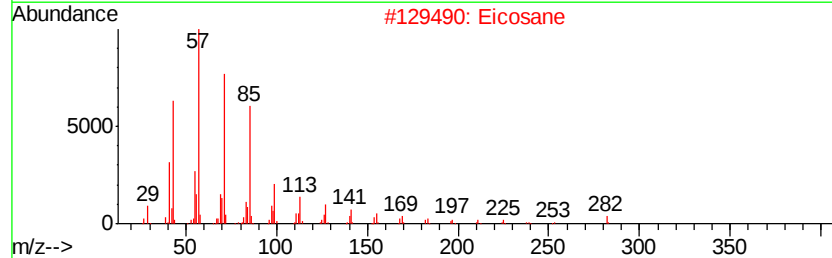
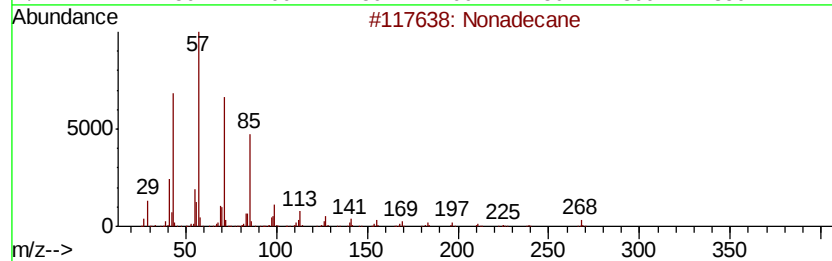
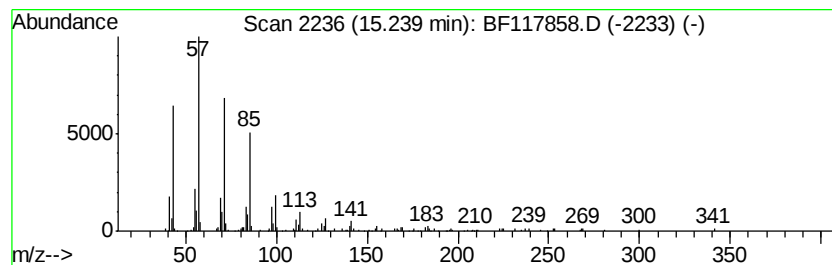
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	2.80 ng	212651	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Eicosane	282	C20H42	000112-95-8	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	87
4		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83
5		Dodecane, 4-methyl-	184	C13H28	006117-97-1	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117858.D
Acq On : 19 Nov 2019 14:42
Operator : JU
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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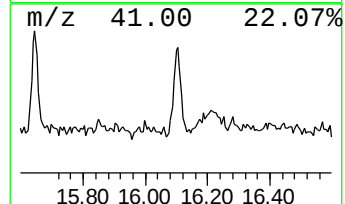
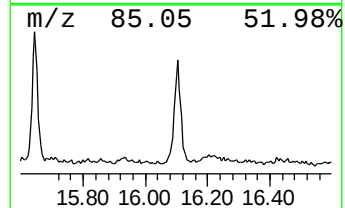
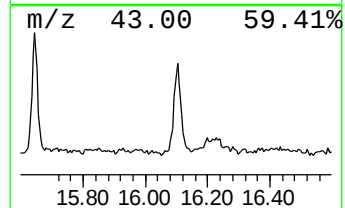
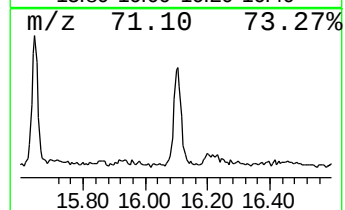
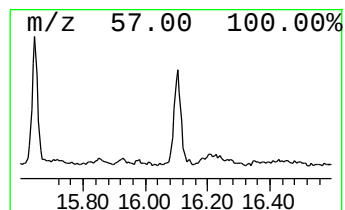
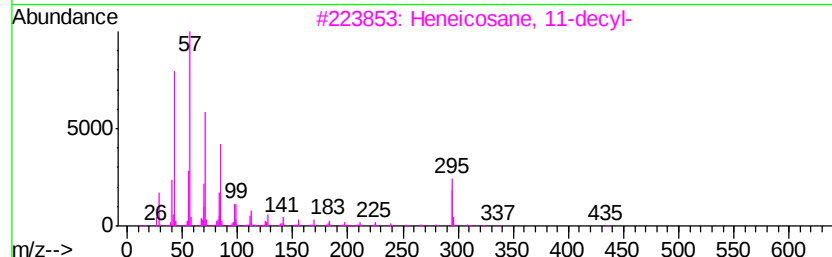
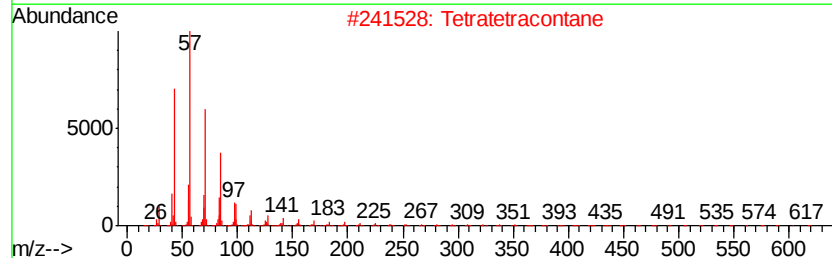
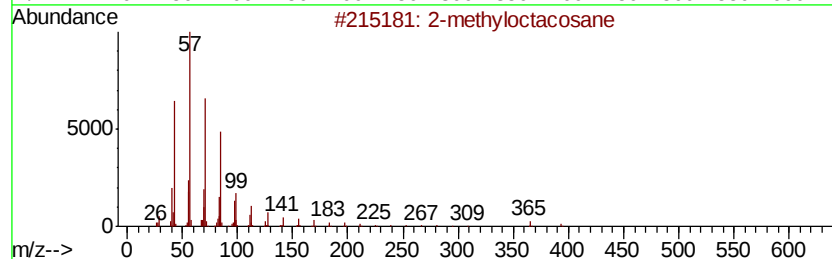
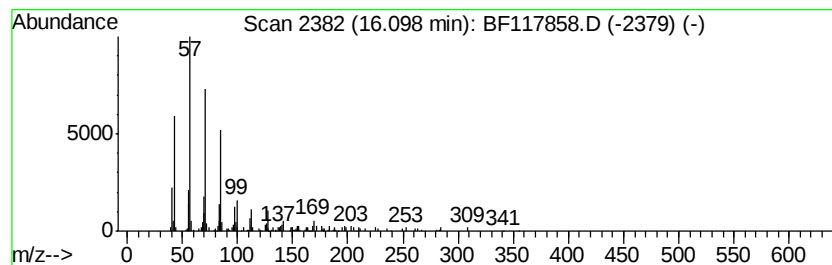
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	2.82 ng	214047	Perylene-d12	15.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-methyloctacosane		408	C29H60	1000376-72-8	90
2	Tetratetracontane		619	C44H90	007098-22-8	90
3	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	87
4	Octacosane		394	C28H58	000630-02-4	87
5	Hexacosane		366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111919\
Data File : BF117858.D
Acq On : 19 Nov 2019 14:42
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Misc :
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ClientSampleId :
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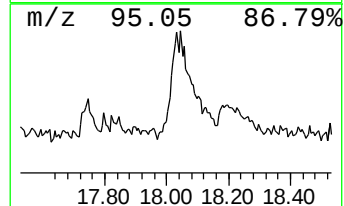
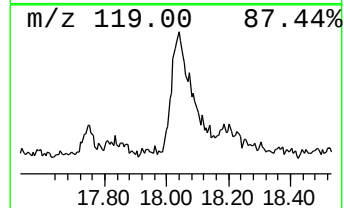
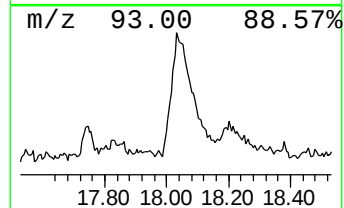
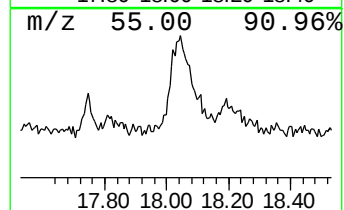
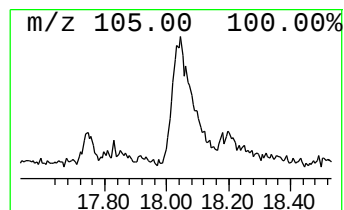
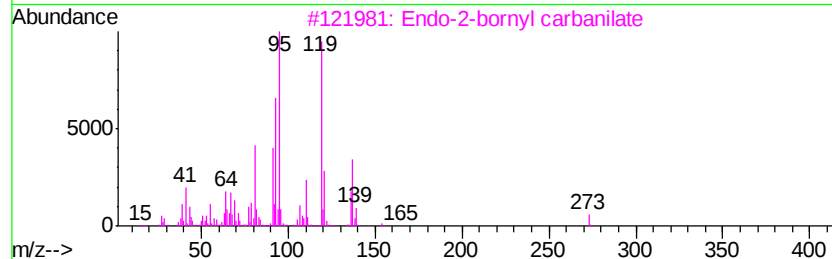
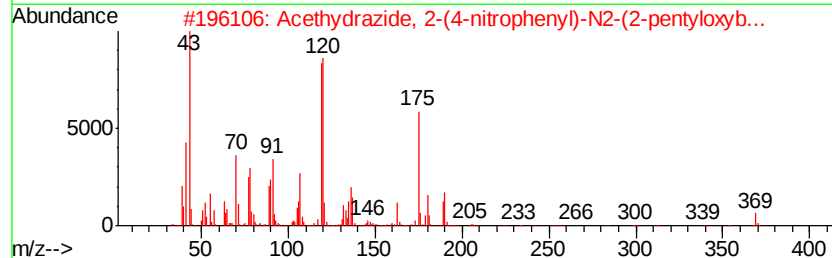
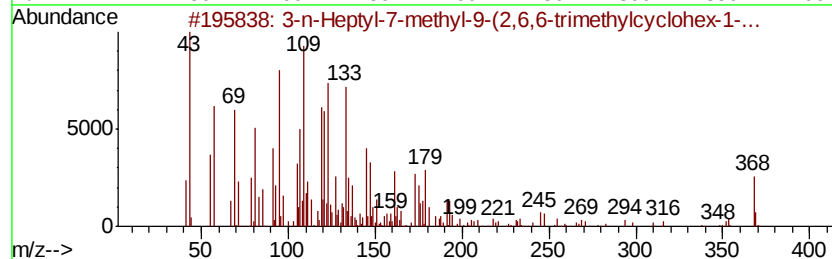
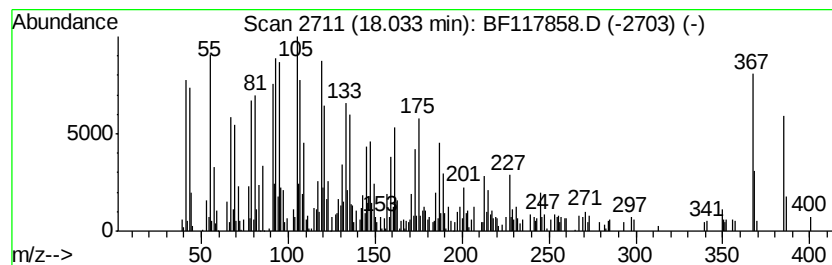
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown18.03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	15.49 ng	1175090	Perylene-d12	15.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-n-Heptyl-7-methyl-9-(2,6,6-tri...	368	C26H40O	1000216-09-3	46
2		Acethydrazide, 2-(4-nitrophenyl)...	369	C20H23N3O4	1000263-97-6	25
3		Endo-2-bornyl carbanilate	273	C17H23NO2	006907-15-9	18
4		Pyrimido[4,5-b]quinoline-2,4(3H,...	368	C17H9ClN4O4	310421-72-8	15
5		2,4-Hexadienamide, N-(4-methylph...	201	C13H15NO	1000362-68-5	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111919\
 Data File : BF117858.D
 Acq On : 19 Nov 2019 14:42
 Operator : JU
 Sample : K5864-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.19	11.7	ng	738606	1	6.92	1259860	20.0
2-Pentanone, 4-hy...	5.13	10.9	ng	689484	1	6.92	1259860	20.0
unknown6.68	6.68	61.6	ng	3880710	1	6.92	1259860	20.0
n-Hexadecanoic acid	11.97	2.5	ng	272087	4	11.46	2156250	20.0
Octadecanoic acid	12.77	4.0	ng	431959	4	11.46	2156250	20.0
Phenol, 4,4'-(1-m...	12.93	3.9	ng	344489	5	14.10	1751100	20.0
Cyclopentadecane	13.96	6.9	ng	604780	5	14.10	1751100	20.0
Tetradecane, 2,6,...	14.89	2.4	ng	182794	6	15.61	1517060	20.0
Nonadecane	15.24	2.8	ng	212651	6	15.61	1517060	20.0
2-methyloctacosane	16.10	2.8	ng	214047	6	15.61	1517060	20.0
unknown18.03	18.03	15.5	ng	1175090	6	15.61	1517060	20.0