

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
 Data File : BF118779.D
 Acq On : 31 Jan 2020 14:01
 Operator : CG/JU
 Sample : PB126458BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB126458BL

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.134	3	8	13	rVB	7349797	11135336	100.00%	13.545%
2	5.051	501	504	511	rVB	97226	122134	1.10%	0.149%
3	5.451	566	572	575	rBV	5059053	5127666	46.05%	6.237%
4	6.451	737	742	750	rBV	3587829	3742890	33.61%	4.553%
5	6.828	802	806	809	rBV	1821349	1634576	14.68%	1.988%
6	6.987	828	833	836	rBV	6508818	6038435	54.23%	7.345%
7	7.387	895	901	904	rBV	4517255	4959711	44.54%	6.033%
8	8.110	1019	1024	1027	rBV	2370187	2176417	19.55%	2.647%
9	9.186	1201	1207	1210	rBV	9012802	9736577	87.44%	11.843%
10	9.575	1269	1273	1276	rBV	766003	597002	5.36%	0.726%
11	9.863	1318	1322	1326	rVB	3163085	2696933	24.22%	3.281%
12	10.651	1450	1456	1460	rBV	6040139	6618778	59.44%	8.051%
13	11.345	1570	1574	1578	rBV	2986925	2660648	23.89%	3.236%
14	11.416	1580	1586	1589	rVB2	360923	310865	2.79%	0.378%
15	11.880	1660	1665	1673	rBV	1371409	1419276	12.75%	1.726%
16	12.663	1794	1798	1811	rBV	678815	756400	6.79%	0.920%
17	12.939	1839	1845	1848	rBV	8848972	9754141	87.60%	11.865%
18	13.833	1993	1997	2004	rBV2	773861	1003826	9.01%	1.221%
19	13.980	2018	2022	2026	rBV	1927107	1815522	16.30%	2.208%
20	14.151	2048	2051	2058	rVB	608909	594332	5.34%	0.723%
21	14.457	2099	2103	2108	rBV	1097057	1074460	9.65%	1.307%
22	14.763	2150	2155	2159	rBV	1359562	1426344	12.81%	1.735%
23	15.098	2207	2212	2222	rVB	1252104	1506193	13.53%	1.832%
24	15.433	2264	2269	2272	rBV	1456023	1809112	16.25%	2.201%
25	15.474	2272	2276	2287	rVB	1038430	1420371	12.76%	1.728%
26	15.904	2343	2349	2354	rBV	749602	1121654	10.07%	1.364%
27	16.398	2427	2433	2441	rVB	372889	706347	6.34%	0.859%
28	16.986	2528	2533	2540	rVB2	125011	244818	2.20%	0.298%

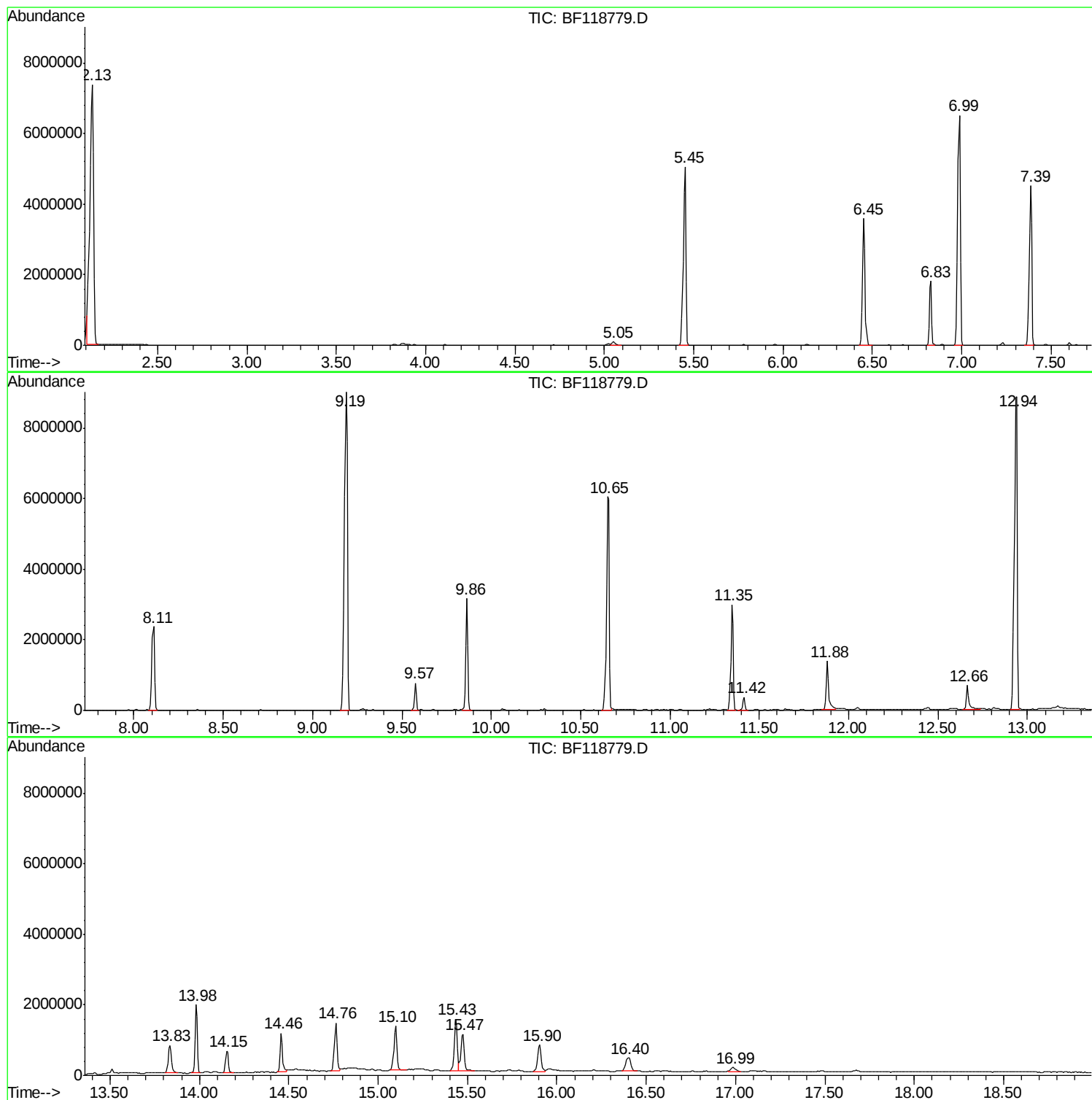
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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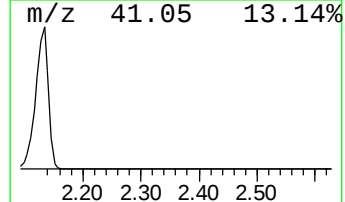
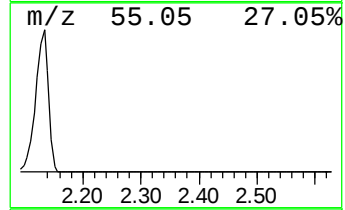
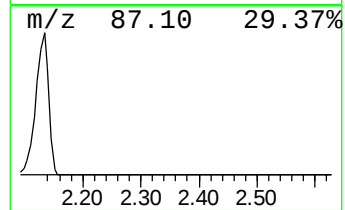
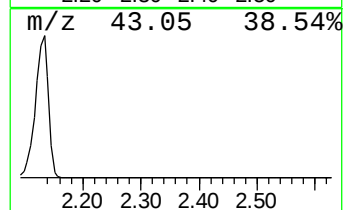
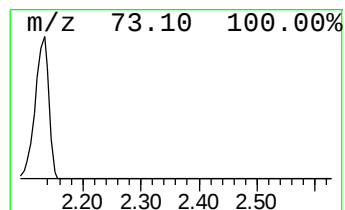
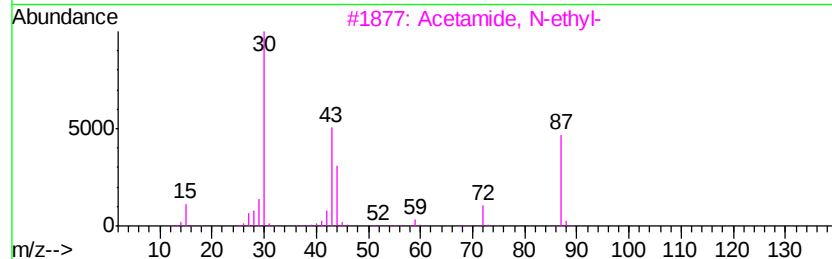
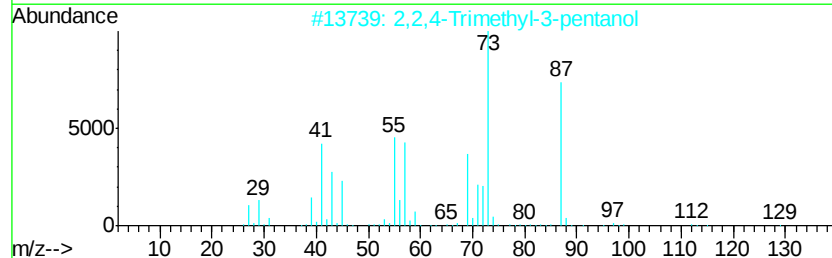
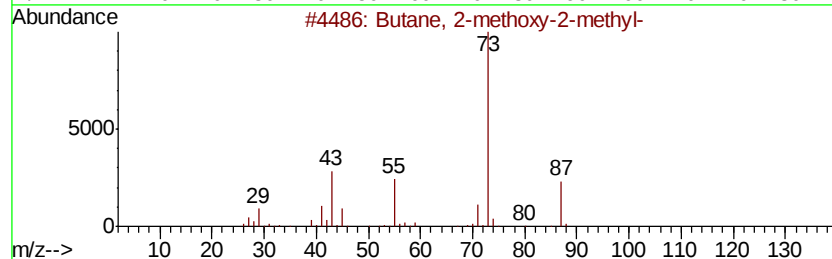
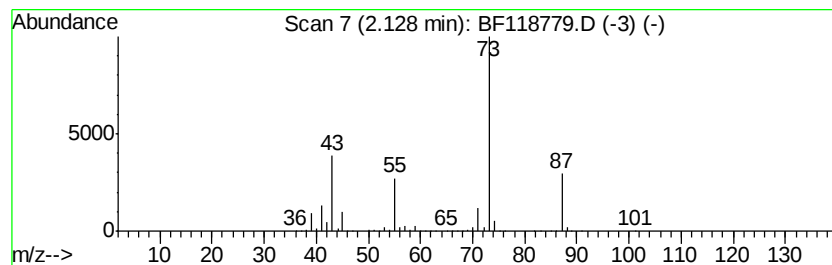
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	136.25 ng	11135300	1,4-Dichlorobenzene-d4	6.83

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	33
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4		1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	17
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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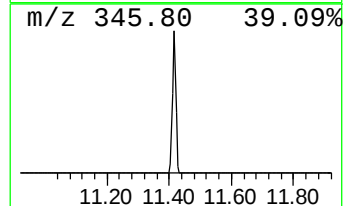
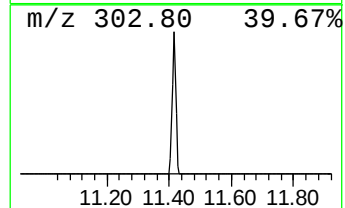
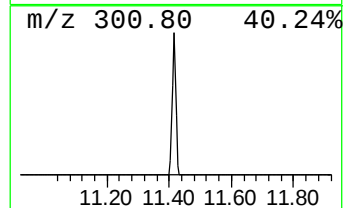
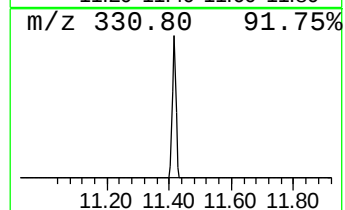
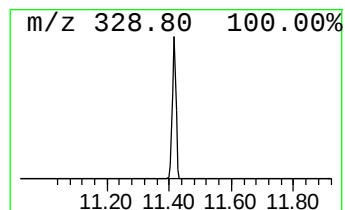
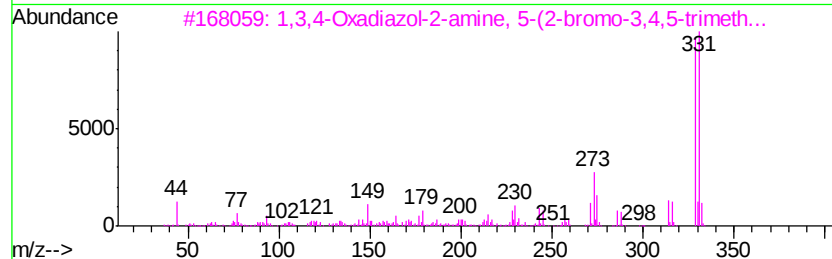
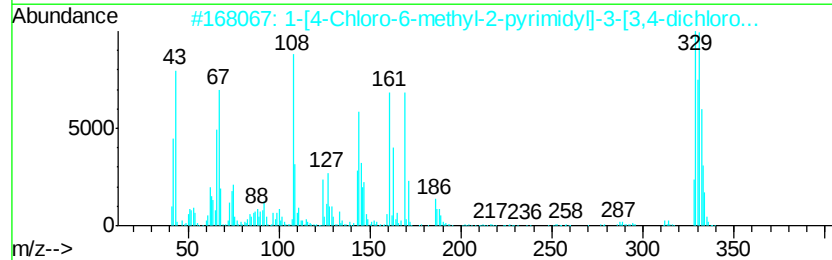
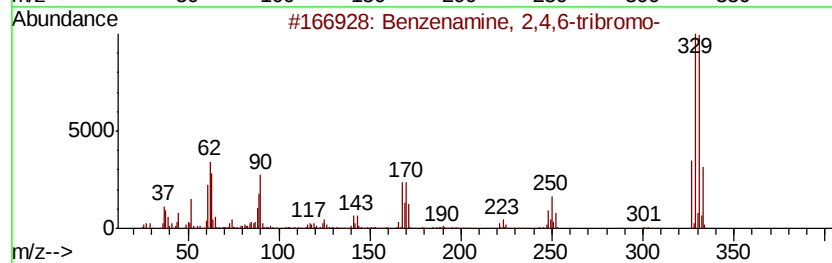
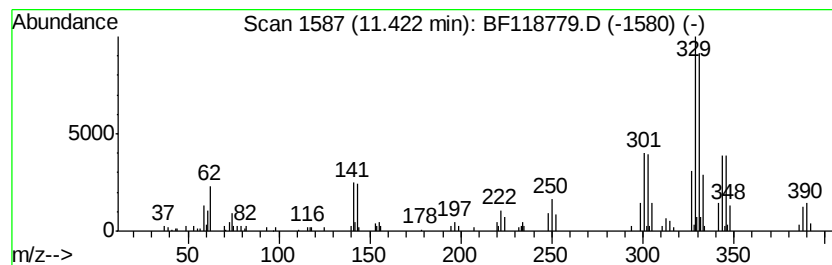
Instrument :
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ClientSampleId :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzenamine, 2,4,6-tribromo- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.42	2.34 ng	310865	Phenanthrene-d10	11.35		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenamine, 2,4,6-tribromo-	327	C6H4Br3N	000147-82-0	86
2		1-[4-Chloro-6-methyl-2-pyrimidyl]...	329	C12H10Cl3N5	051387-79-2	11
3		1,3,4-Oxadiazol-2-amine, 5-(2-br...	329	C11H12BrN3O4	1000350-74-8	10
4		6-(4-Chlorophenyl)benzo[4,5]imid...	329	C20H12ClN3	105997-06-6	9
5		4,4'-(p-Phenylenediisopropyliden...	344	C24H28N2	002716-10-1	9



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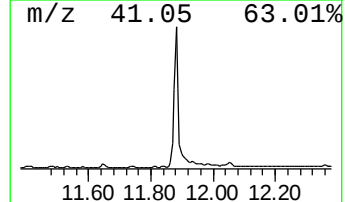
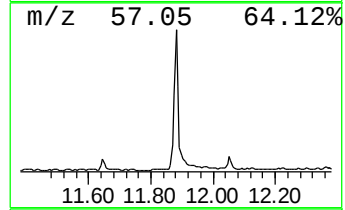
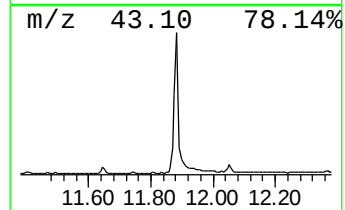
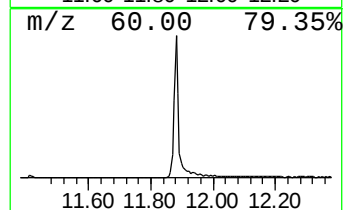
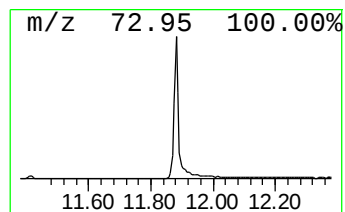
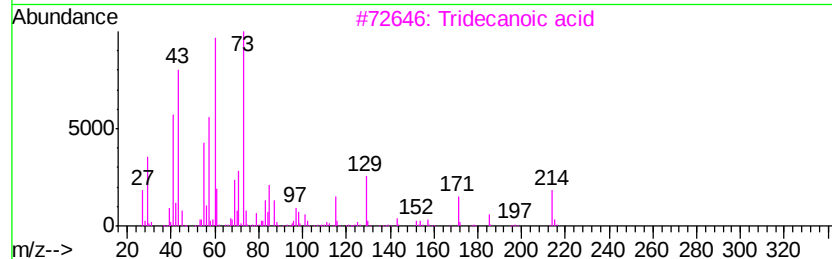
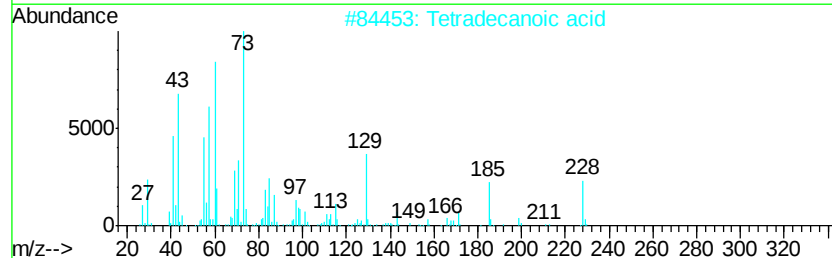
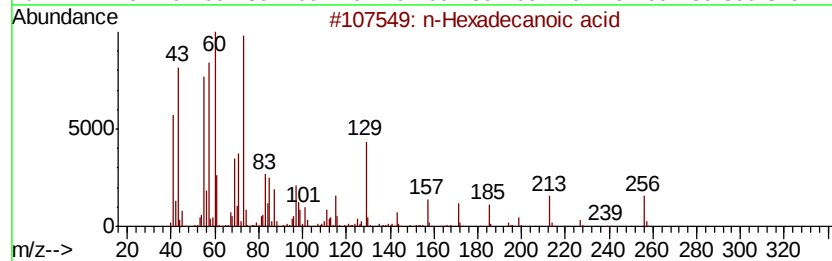
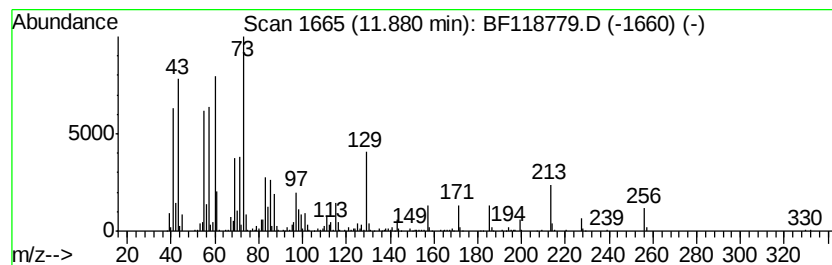
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	10.67 ng	1419280	Phenanthrene-d10	11.35

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5		Octadecanoic acid	284	C18H36O2	000057-11-4	81



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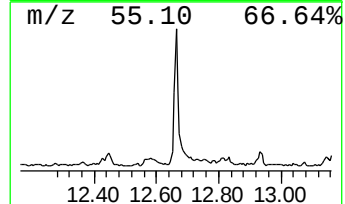
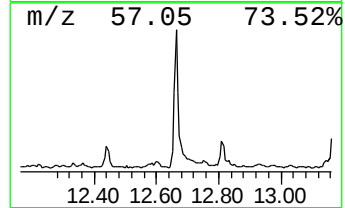
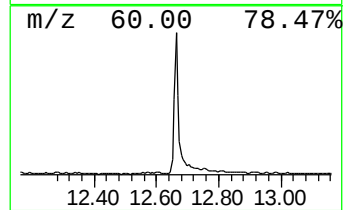
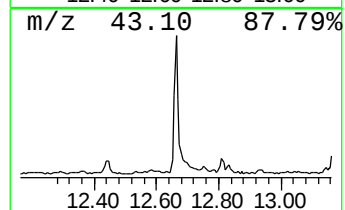
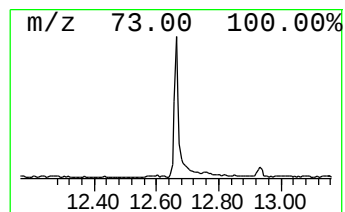
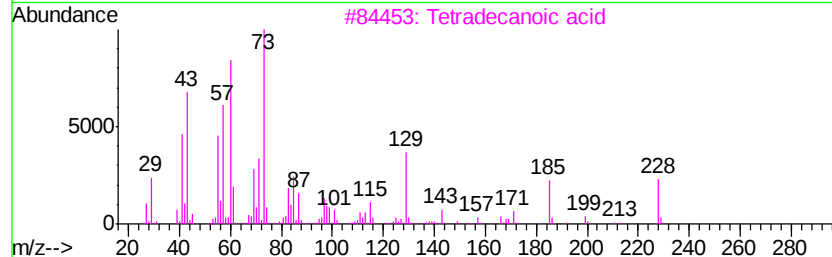
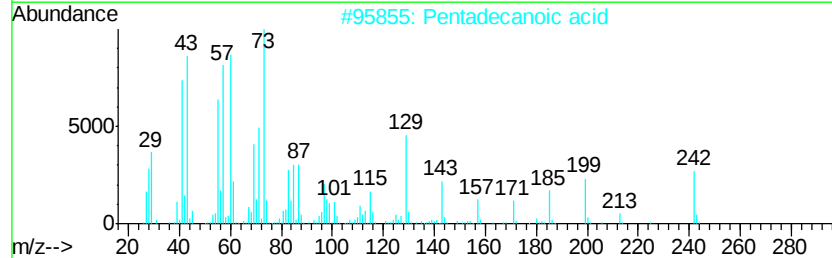
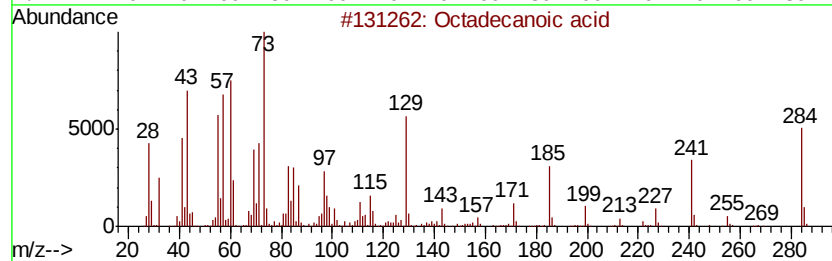
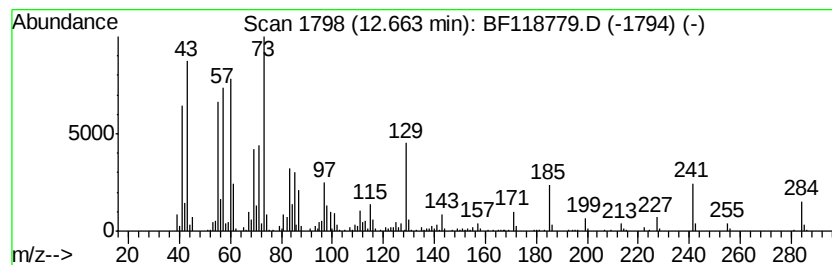
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Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	5.69 ng	756400	Phenanthrene-d10	11.35

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	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
5		Tridecanoic acid	214	C13H26O2	000638-53-9	74



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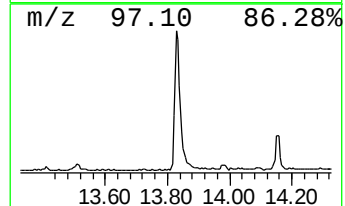
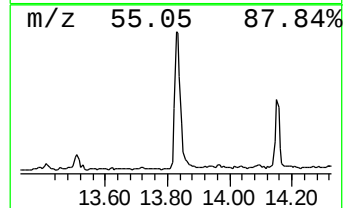
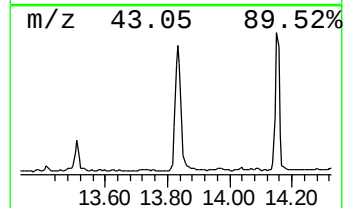
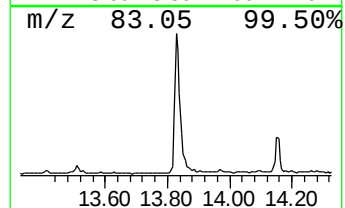
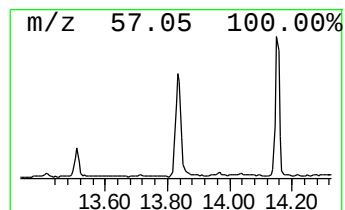
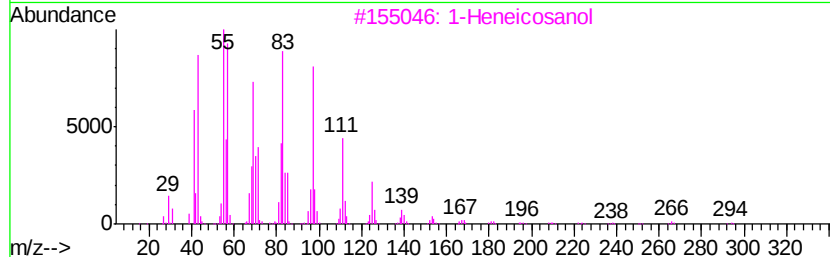
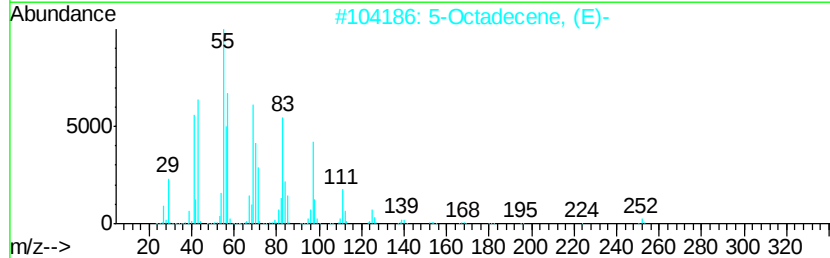
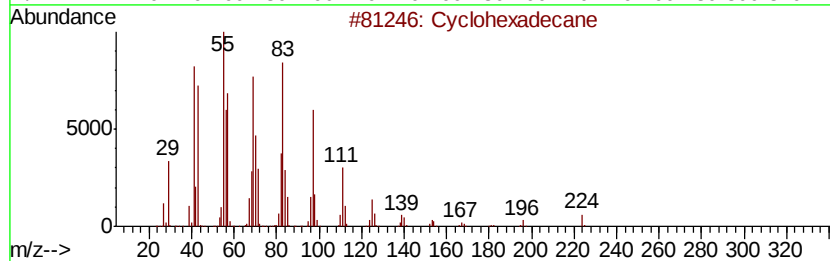
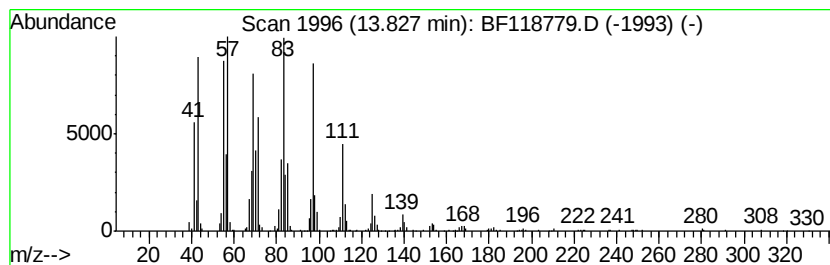
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclohexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	11.06 ng	1003830	Chrysene-d12	13.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	99
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	93
3		1-Heneicosanol	312	C21H44O	015594-90-8	93
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5		Behenic alcohol	326	C22H46O	000661-19-8	93



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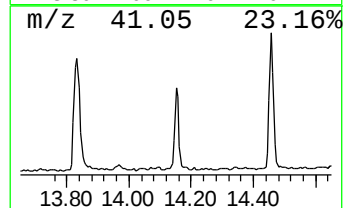
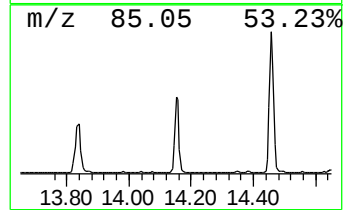
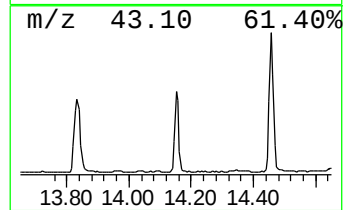
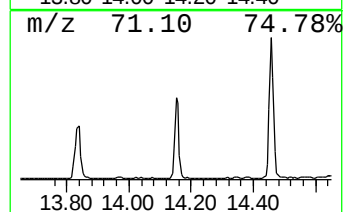
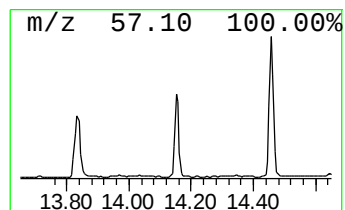
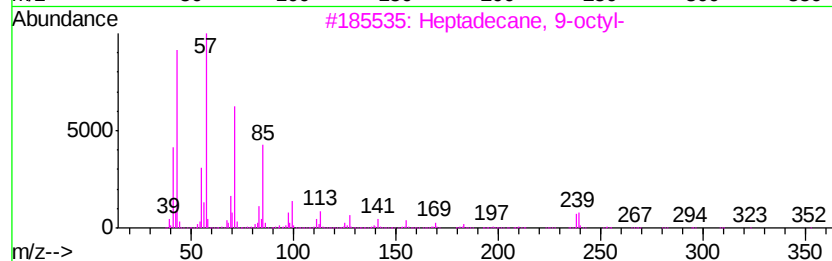
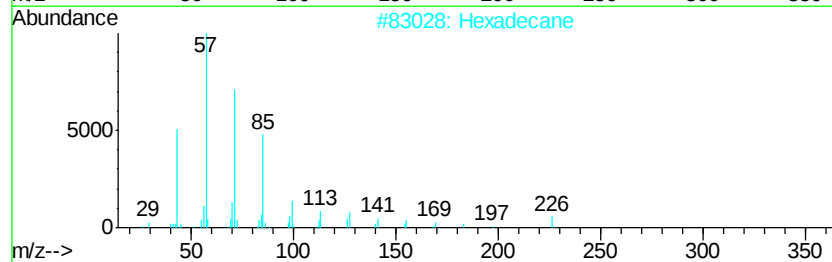
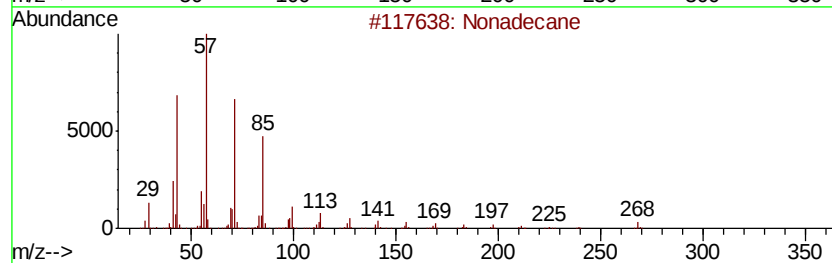
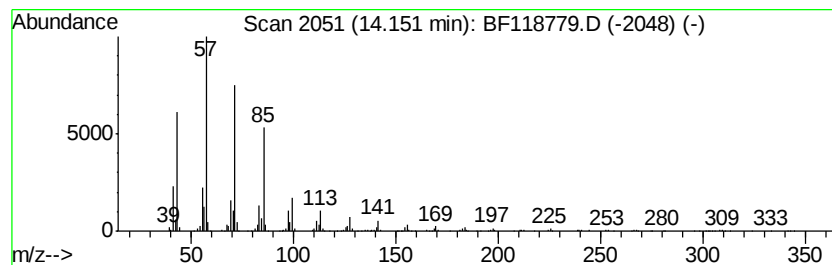
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	6.55 ng	594332	Chrysene-d12	13.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Hexadecane	226	C16H34	000544-76-3	93
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
Data File : BF118779.D
Acq On : 31 Jan 2020 14:01
Operator : CG/JU
Sample : PB126458BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB126458BL

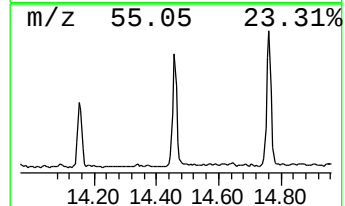
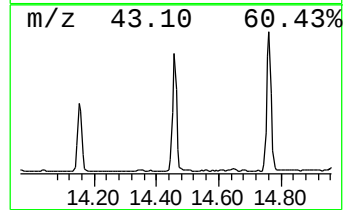
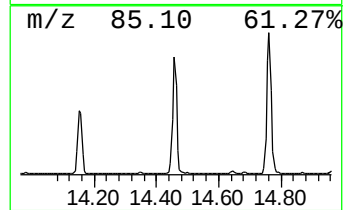
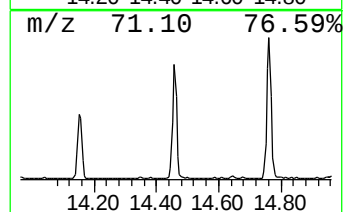
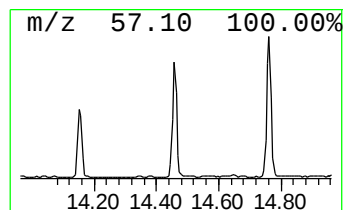
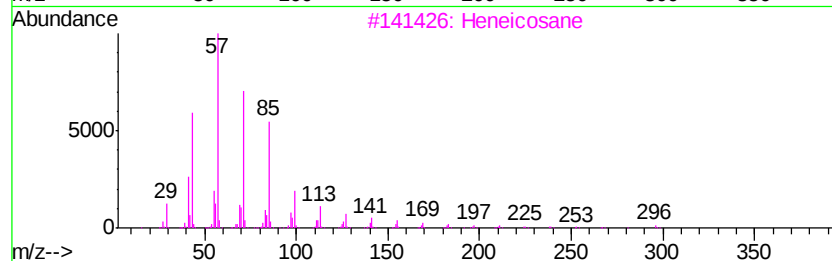
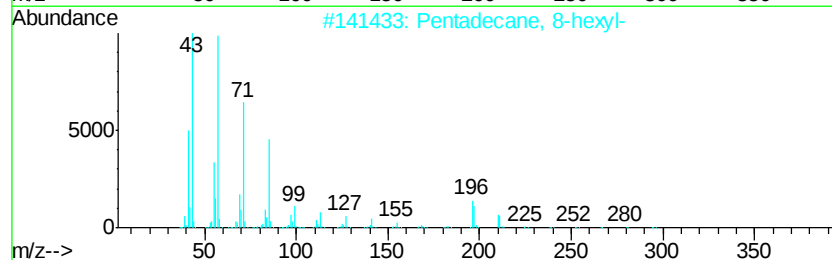
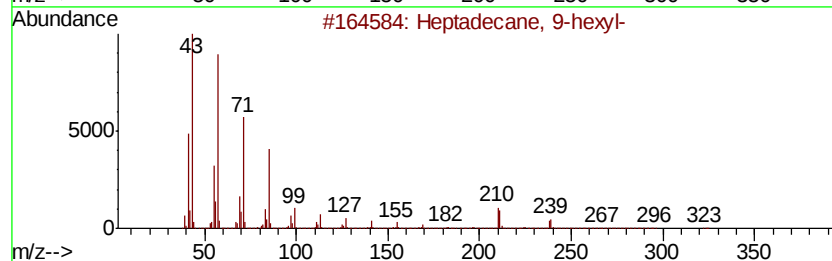
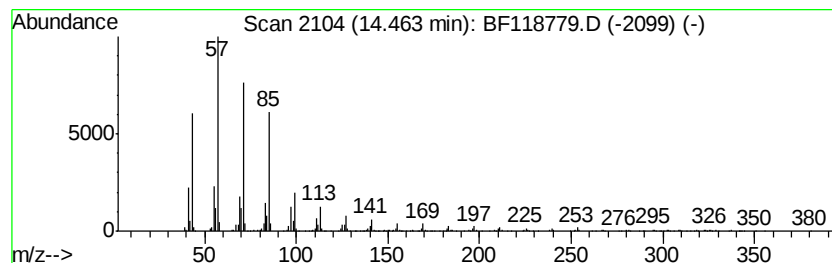
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptadecane, 9-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	11.84 ng	1074460	Chrysene-d12	13.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
3		Heneicosane	296	C21H44	000629-94-7	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
Data File : BF118779.D
Acq On : 31 Jan 2020 14:01
Operator : CG/JU
Sample : PB126458BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

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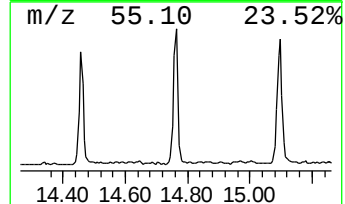
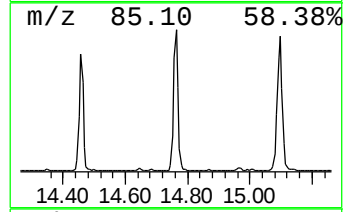
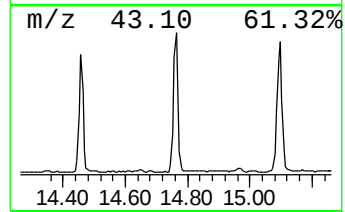
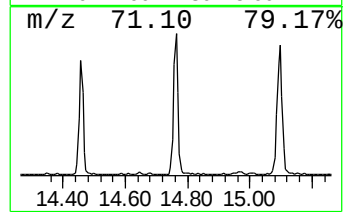
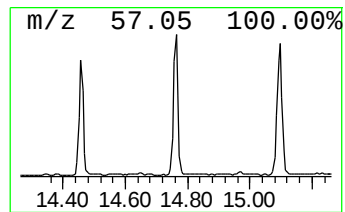
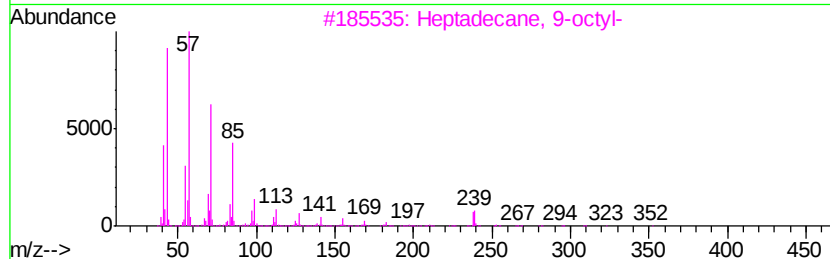
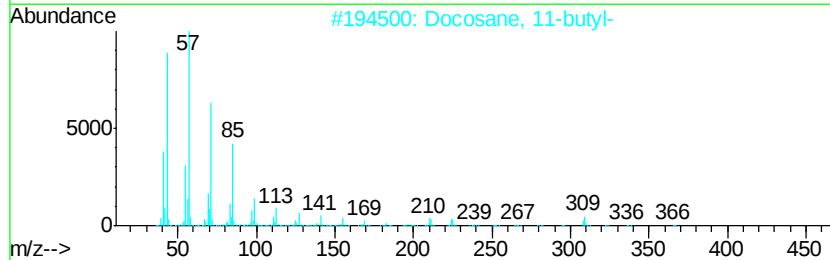
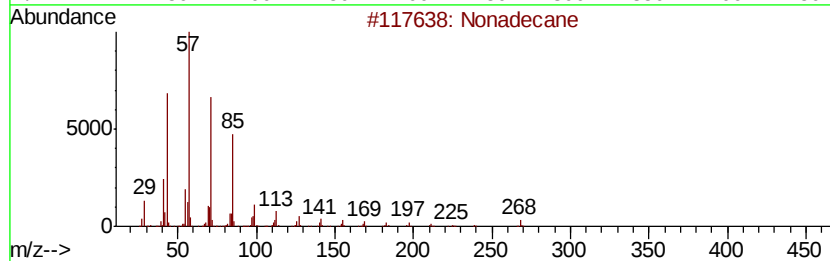
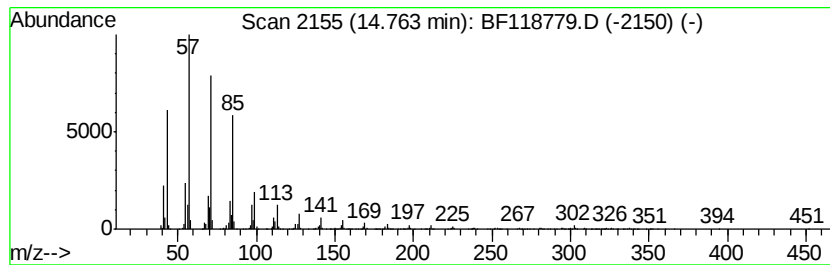
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Docosane, 11-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	15.77 ng	1426340	Perylene-d12	15.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	94
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4		Hexadecane	226	C16H34	000544-76-3	93
5		Eicosane, 9-octyl-	394	C28H58	013475-77-9	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
Data File : BF118779.D
Acq On : 31 Jan 2020 14:01
Operator : CG/JU
Sample : PB126458BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

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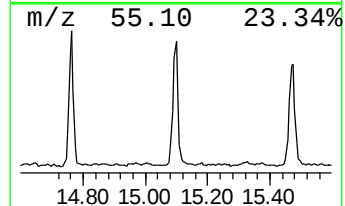
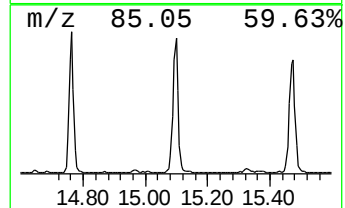
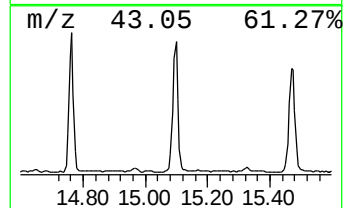
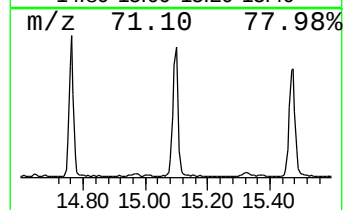
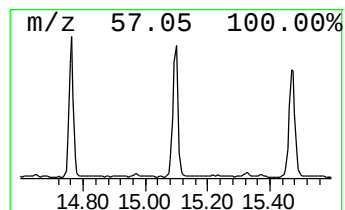
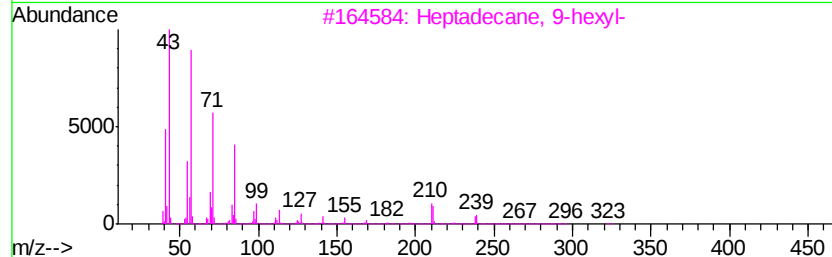
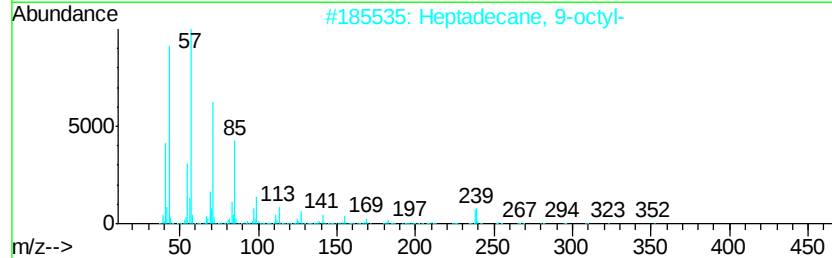
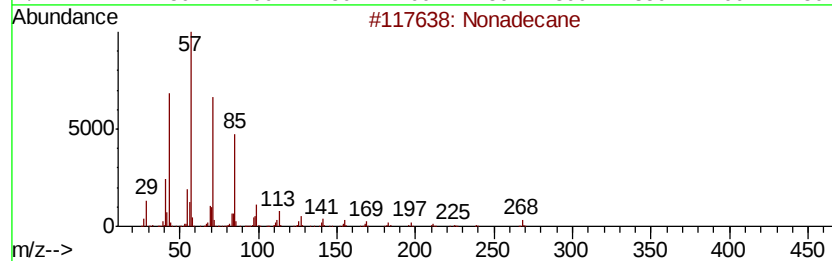
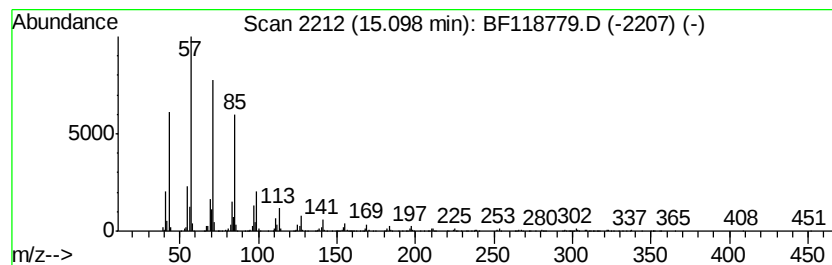
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	16.65 ng	1506190	Perylene-d12	15.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	93
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
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 Acq On : 31 Jan 2020 14:01
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 Misc :
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 ClientSampleId :
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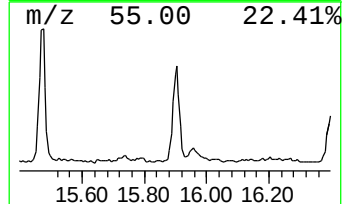
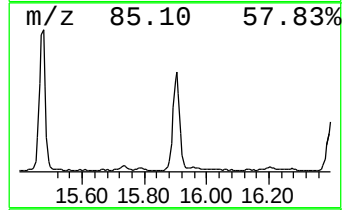
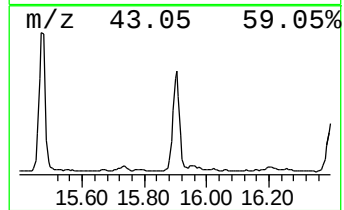
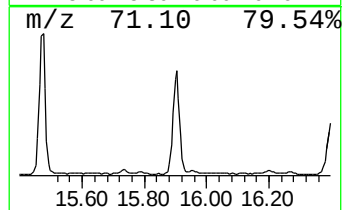
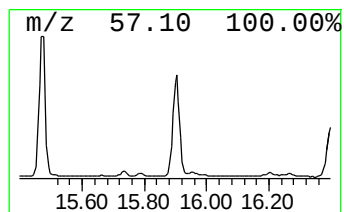
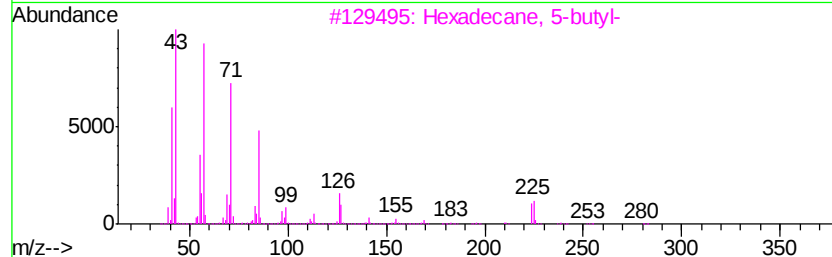
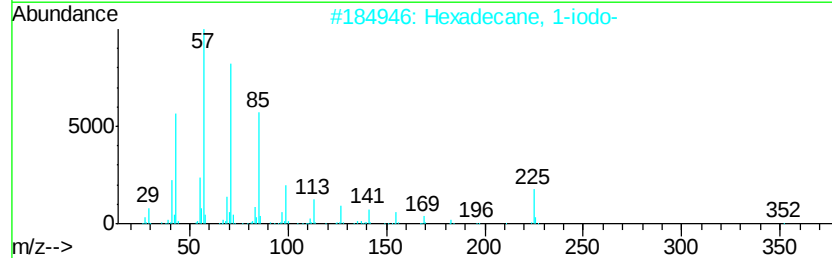
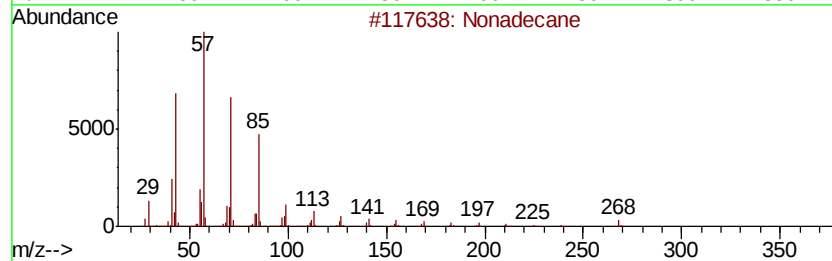
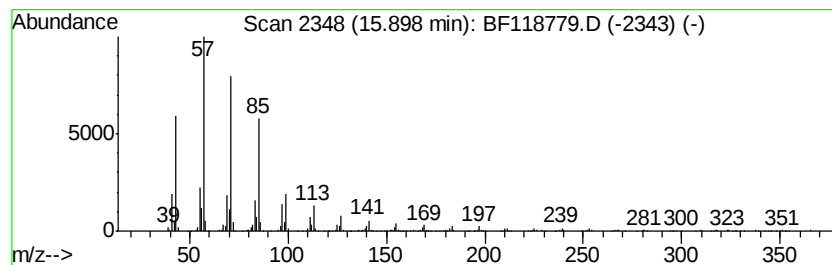
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Hexadecane, 1-iodo- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	12.40 ng	1121650	Perylene-d12	15.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	94
3		Hexadecane, 5-butyl-	282	C20H42	006912-07-8	93
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\BF013120\
Data File : BF118779.D
Acq On : 31 Jan 2020 14:01
Operator : CG/JU
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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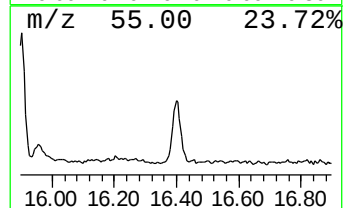
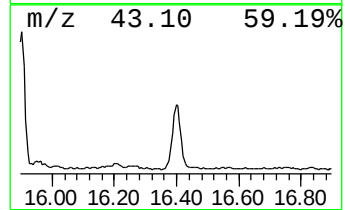
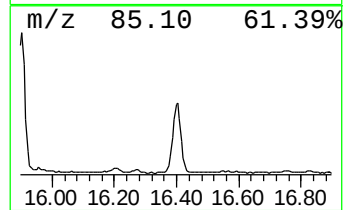
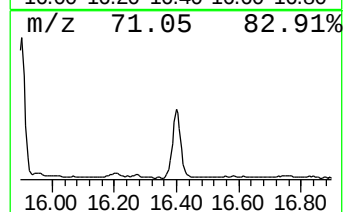
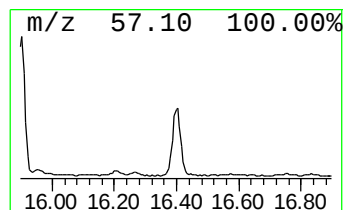
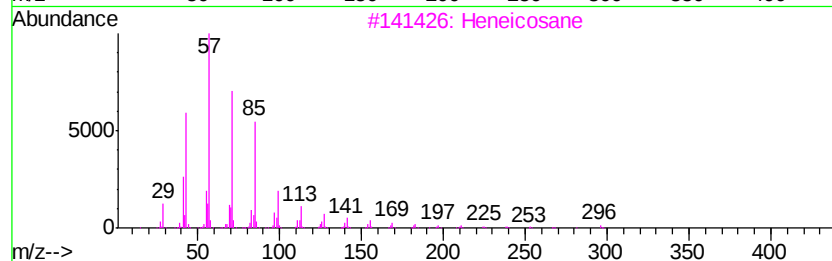
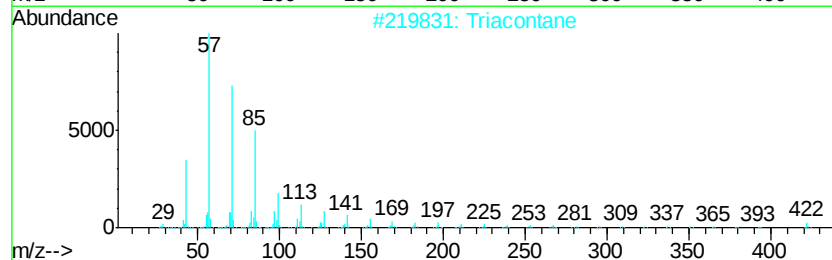
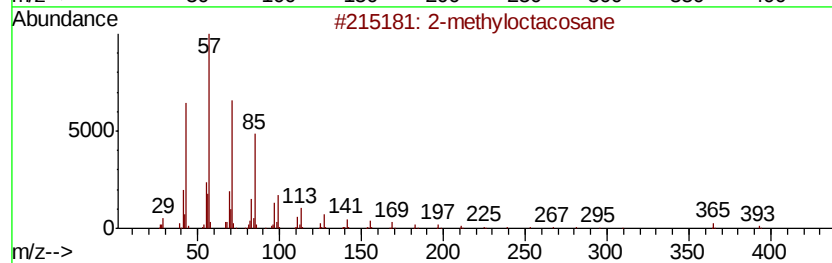
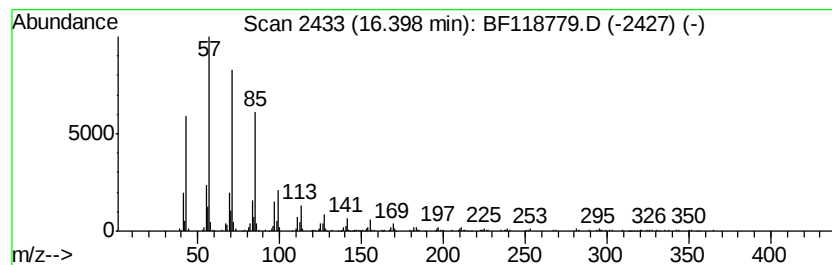
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	7.81 ng	706347	Perylene-d12	15.43

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Triacotane	422	C30H62	000638-68-6	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			2-methylhexacosane	380	C27H56	1000376-72-7	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
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Acq On : 31 Jan 2020 14:01
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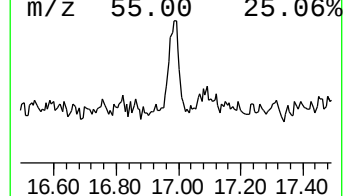
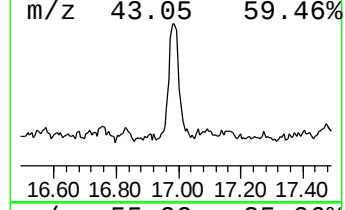
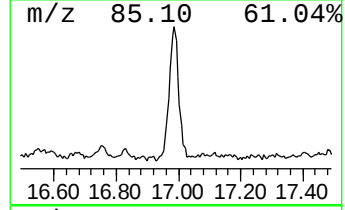
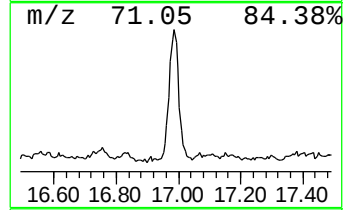
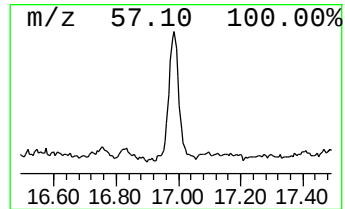
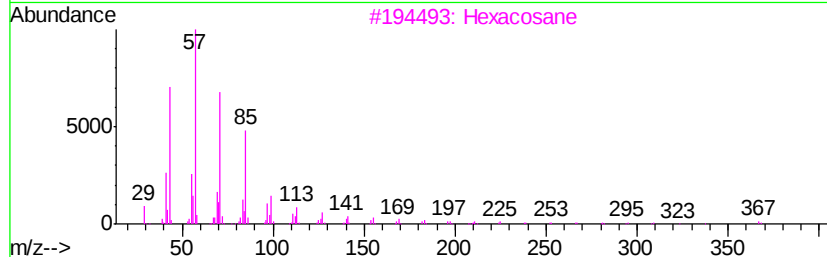
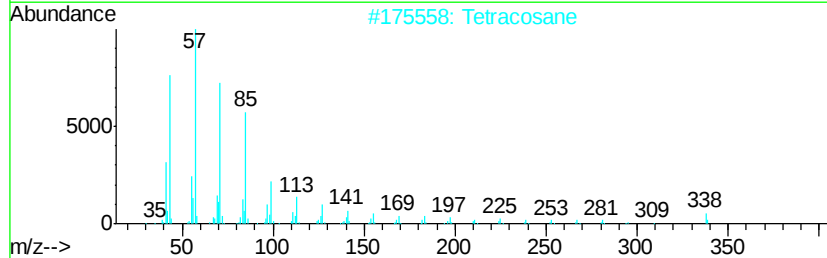
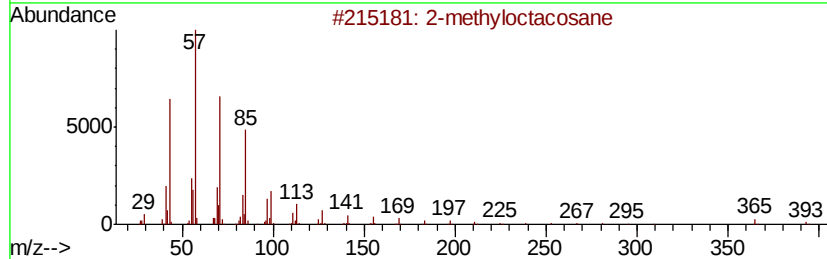
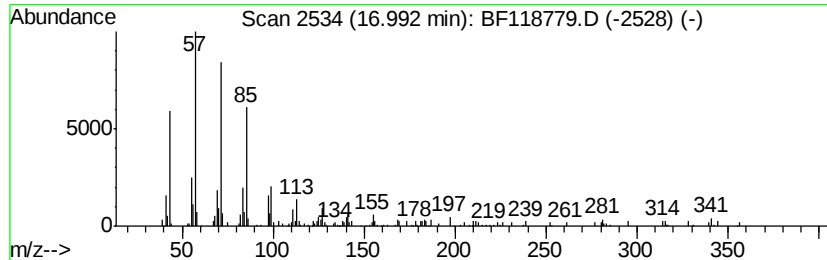
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.99	2.71 ng	244818	Perylene-d12	15.43

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Tetracosane	338	C24H50	000646-31-1	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87
5			Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF013120\
 Data File : BF118779.D
 Acq On : 31 Jan 2020 14:01
 Operator : CG/JU
 Sample : PB126458BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB126458BL

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...	2.13	136.3	ng	11135300	1	6.83	1634580	20.0
Benzenamine, 2,4,...	11.42	2.3	ng	310865	4	11.35	2660650	20.0
n-Hexadecanoic acid	11.88	10.7	ng	1419280	4	11.35	2660650	20.0
Octadecanoic acid	12.66	5.7	ng	756400	4	11.35	2660650	20.0
Cyclohexadecane	13.83	11.1	ng	1003830	5	13.98	1815520	20.0
Nonadecane	14.15	6.5	ng	594332	5	13.98	1815520	20.0
Heptadecane, 9-he...	14.46	11.8	ng	1074460	5	13.98	1815520	20.0
Docosane, 11-butyl-	14.76	15.8	ng	1426340	6	15.43	1809110	20.0
Heptadecane, 9-oc...	15.10	16.6	ng	1506190	6	15.43	1809110	20.0
Hexadecane, 1-iodo-	15.90	12.4	ng	1121650	6	15.43	1809110	20.0
2-methyloctacosane	16.40	7.8	ng	706347	6	15.43	1809110	20.0
Tetracosane	16.99	2.7	ng	244818	6	15.43	1809110	20.0