

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
 Data File : BP002761.D
 Acq On : 15 Jul 2020 19:03
 Operator : CG/JU
 Sample : L3304-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-GF-02-058-B

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_P\METHODS\8270-BP071020.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	4.716	304	311	323	rBV	68639	103733	2.56%	0.220%
2	5.187	384	391	412	rBV	1688495	2682999	66.22%	5.679%
3	6.763	652	659	671	rBV	1616111	2820722	69.62%	5.971%
4	6.863	671	676	687	rVB	48371	97234	2.40%	0.206%
5	7.210	727	735	743	rBV4	21679	57744	1.43%	0.122%
6	7.558	787	794	804	rBV	681217	1052824	25.99%	2.229%
7	8.704	972	989	1001	rBV	1041946	1862617	45.97%	3.943%
8	8.805	1002	1006	1015	rVB2	36923	66188	1.63%	0.140%
9	10.328	1256	1265	1272	rBV	854395	1517663	37.46%	3.212%
10	11.393	1439	1446	1451	rBV3	42261	89770	2.22%	0.190%
11	11.793	1506	1514	1521	rBV	71436	149528	3.69%	0.317%
12	12.010	1541	1551	1559	rVB4	40221	113128	2.79%	0.239%
13	12.228	1583	1588	1591	rBV	27985	43062	1.06%	0.091%
14	12.269	1591	1595	1600	rVB3	24581	41997	1.04%	0.089%
15	12.628	1652	1656	1663	rBV5	22837	44233	1.09%	0.094%
16	12.763	1675	1679	1683	rBV2	43766	61810	1.53%	0.131%
17	12.822	1683	1689	1700	rVV	2368718	3617531	89.29%	7.657%
18	13.057	1725	1729	1736	rVB	91321	130636	3.22%	0.277%
19	13.369	1777	1782	1790	rBV3	45929	108490	2.68%	0.230%
20	13.534	1805	1810	1814	rBV2	66592	100538	2.48%	0.213%
21	13.581	1814	1818	1823	rVB3	50402	78713	1.94%	0.167%
22	13.634	1823	1827	1829	rBV	45624	61089	1.51%	0.129%
23	13.669	1829	1833	1838	rVV	298784	417651	10.31%	0.884%
24	13.728	1839	1843	1846	rVB3	49954	63983	1.58%	0.135%
25	13.928	1873	1877	1885	rBV2	113879	175667	4.34%	0.372%
26	14.145	1909	1914	1918	rBV	134105	169627	4.19%	0.359%
27	14.210	1918	1925	1931	rVV	1188810	1813008	44.75%	3.838%
28	14.269	1931	1935	1940	rVV	82015	118766	2.93%	0.251%
29	14.428	1958	1962	1966	rBV5	24782	51412	1.27%	0.109%
30	14.616	1990	1994	1999	rVB3	63289	98195	2.42%	0.208%
31	14.698	2005	2008	2013	rVB2	44782	60595	1.50%	0.128%
32	14.769	2016	2020	2025	rVB2	60059	94109	2.32%	0.199%
33	14.839	2028	2032	2036	rBV2	53006	86989	2.15%	0.184%
34	15.010	2058	2061	2063	rBV2	35572	40668	1.00%	0.086%

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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_P\METHODS\8270-BP071020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.039	2063	2066	2072	rVB2	69316	82541	2.04%	0.175%
36	15.104	2073	2077	2082	rBV	150126	190267	4.70%	0.403%
37	15.269	2100	2105	2109	rBV	118892	178674	4.41%	0.378%
38	15.516	2144	2147	2151	rVB	92883	119525	2.95%	0.253%
39	15.710	2175	2180	2189	rVB	1483058	2196727	54.22%	4.650%
40	15.969	2220	2224	2226	rBV	187165	233073	5.75%	0.493%
41	15.992	2226	2228	2236	rVB	165552	223092	5.51%	0.472%
42	16.381	2291	2294	2298	rVB4	38331	46550	1.15%	0.099%
43	16.763	2356	2359	2365	rBV2	158831	270345	6.67%	0.572%
44	16.816	2365	2368	2375	rVB	154190	236977	5.85%	0.502%
45	16.969	2389	2394	2398	rBV	1323642	1912982	47.22%	4.049%
46	17.010	2398	2401	2406	rVB	896189	1164713	28.75%	2.465%
47	17.098	2413	2416	2421	rVB	258218	371193	9.16%	0.786%
48	17.504	2482	2485	2489	rVB2	161711	187187	4.62%	0.396%
49	17.675	2510	2514	2523	rVB	643128	791265	19.53%	1.675%
50	17.851	2540	2544	2547	rBV	133511	175865	4.34%	0.372%
51	17.892	2548	2551	2558	rVB	243240	320888	7.92%	0.679%
52	18.033	2571	2575	2579	rBV2	224891	358293	8.84%	0.758%
53	18.186	2598	2601	2605	rVB	171991	182256	4.50%	0.386%
54	18.339	2624	2627	2630	rBV2	94302	114863	2.84%	0.243%
55	18.633	2674	2677	2680	rBV3	129601	155243	3.83%	0.329%
56	18.745	2693	2696	2699	rBV	138206	168935	4.17%	0.358%
57	18.786	2699	2703	2705	rBV	1837864	2145850	52.96%	4.542%
58	18.816	2705	2708	2713	rVB	3367938	4051514	100.00%	8.576%
59	18.951	2727	2731	2734	rBV	359581	417960	10.32%	0.885%
60	18.998	2734	2739	2746	rVB	1371789	1765531	43.58%	3.737%
61	19.351	2792	2799	2809	rVB	1323716	2174408	53.67%	4.603%
62	19.557	2830	2834	2839	rVB2	3191800	3853666	95.12%	8.157%
63	19.886	2887	2890	2892	rBV	181180	200619	4.95%	0.425%
64	20.822	3046	3049	3053	rVB	452265	516050	12.74%	1.092%
65	21.074	3087	3092	3096	rBV2	1476292	2517017	62.13%	5.328%
66	21.116	3096	3099	3103	rVB	535818	624803	15.42%	1.323%
67	23.274	3462	3466	3471	rVB2	711952	1200943	29.64%	2.542%

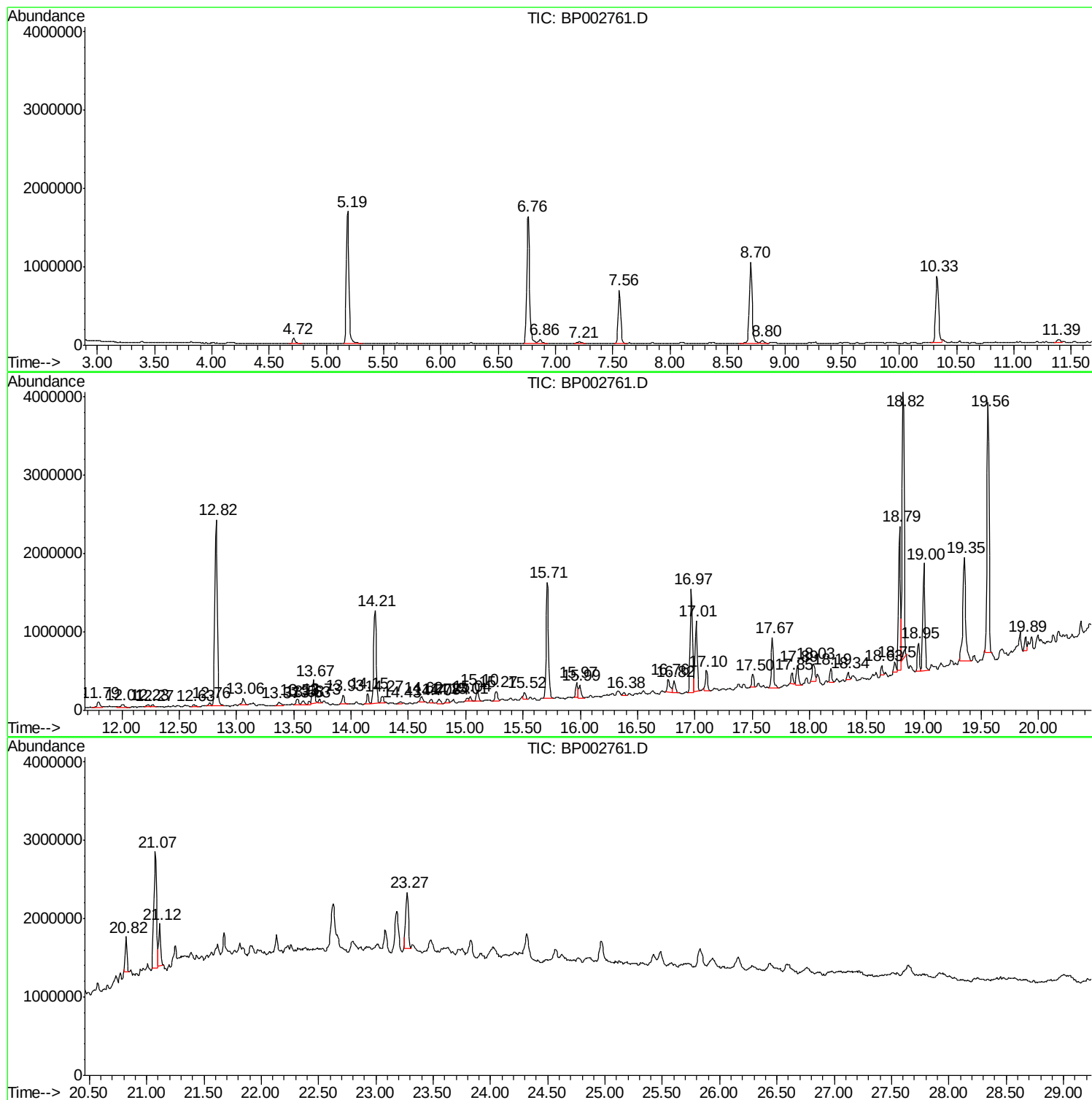
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Data File : BP002761.D
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Sample : L3304-03
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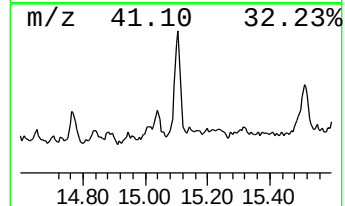
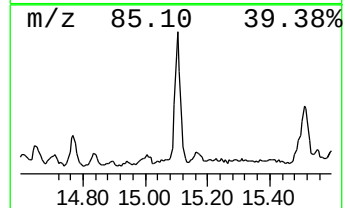
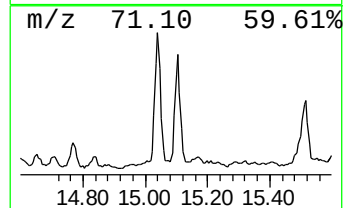
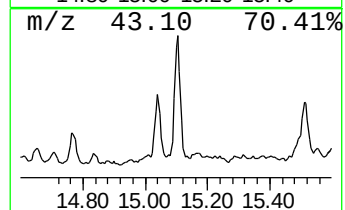
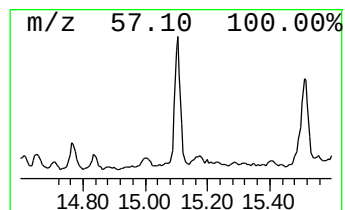
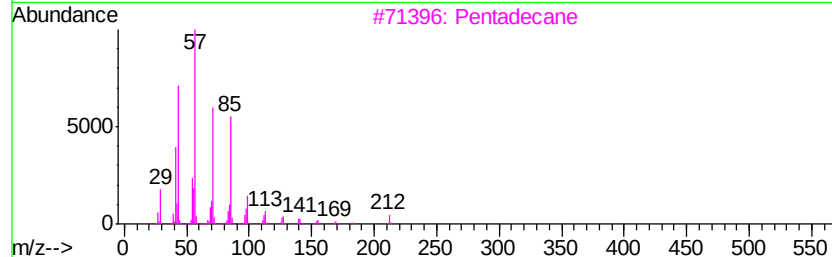
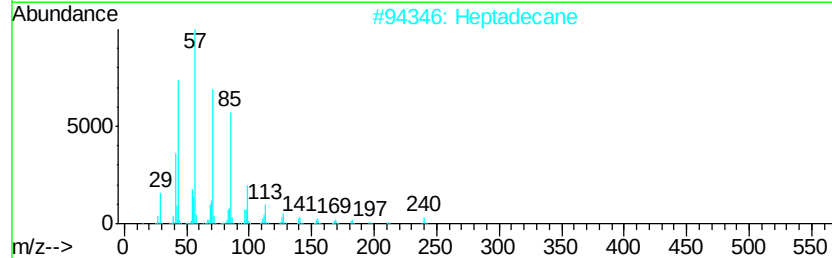
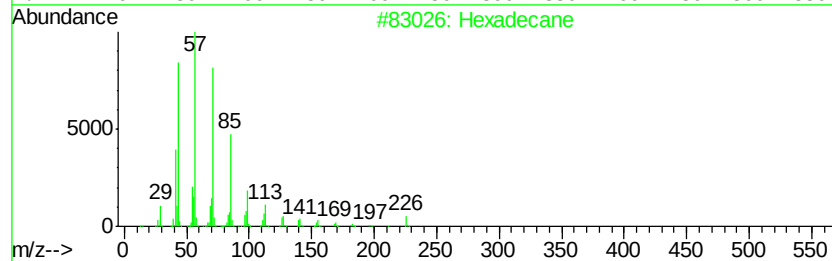
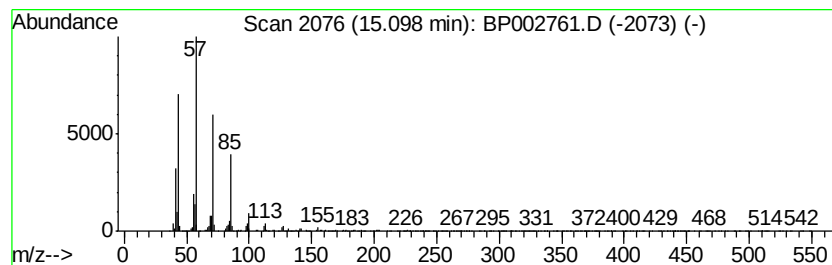
Instrument :
BNA_P
ClientSampleId :
FK-GF-02-058-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.10	2.10 ng	190267	Acenaphthene-d10		14.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Heptadecane	240	C17H36	000629-78-7	90
3	Pentadecane	212	C15H32	000629-62-9	90
4	Heneicosane	296	C21H44	000629-94-7	81
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	80



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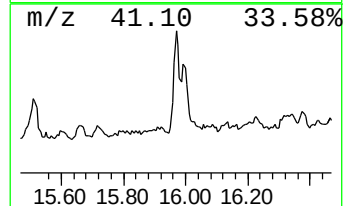
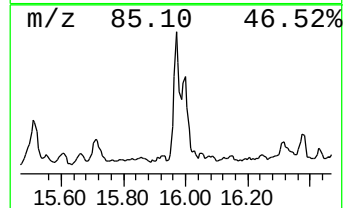
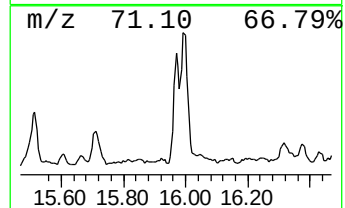
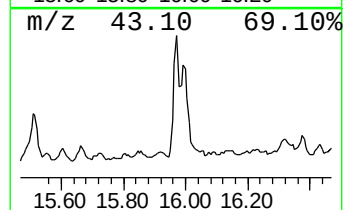
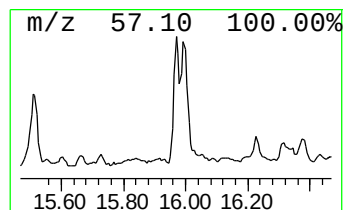
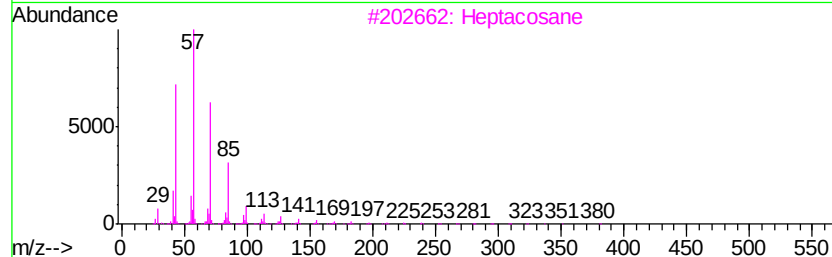
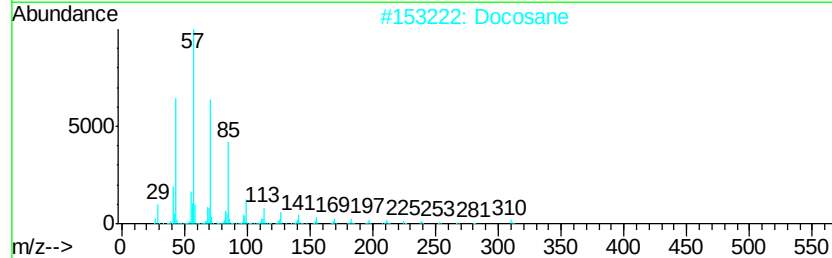
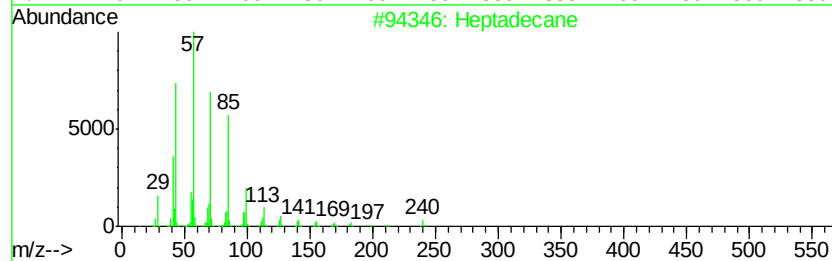
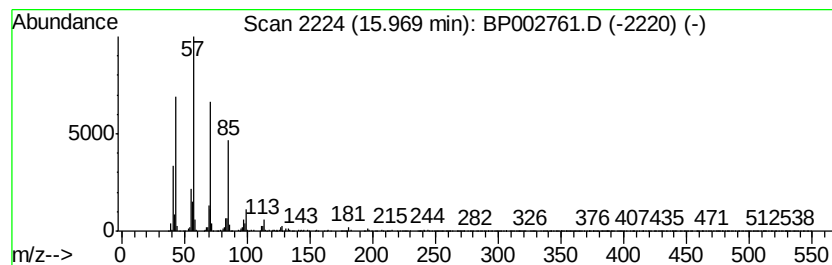
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Heptadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	2.44 ng	233073	Phenanthrene-d10	16.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	97
2	Docosane		310	C22H46	000629-97-0	91
3	Heptacosane		380	C27H56	000593-49-7	91
4	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
5	Tetracosane		338	C24H50	000646-31-1	90



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Misc :
ALS Vial : 13 Sample Multiplier: 1

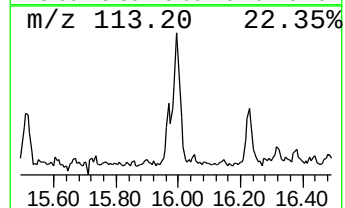
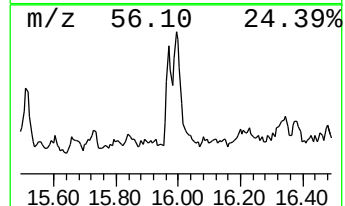
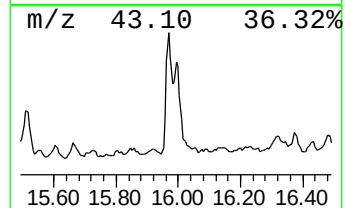
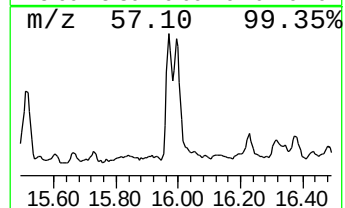
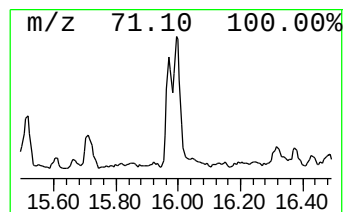
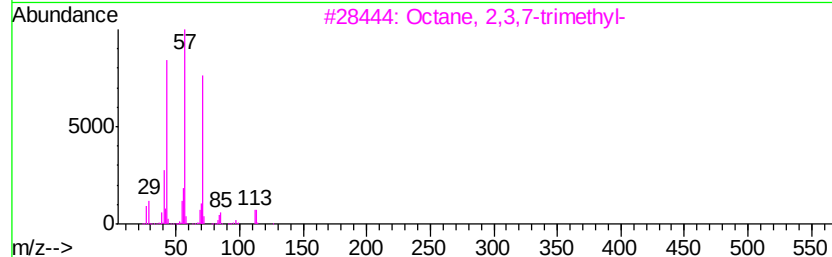
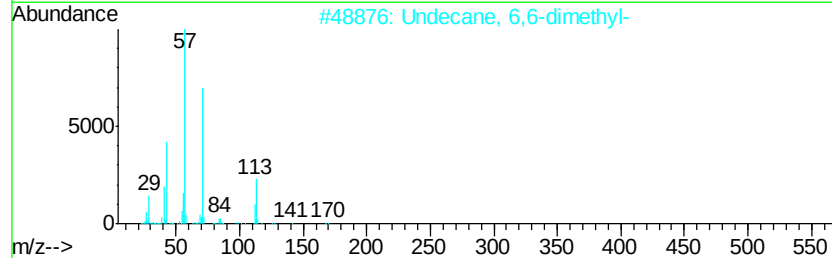
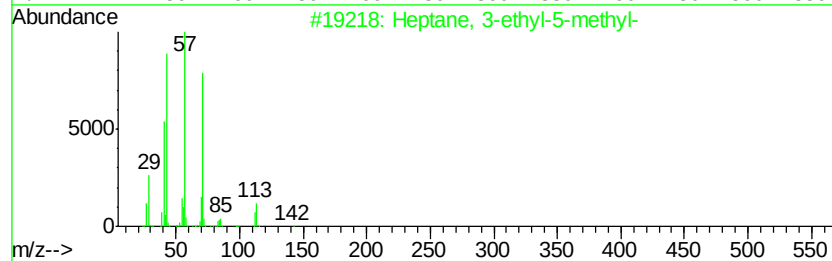
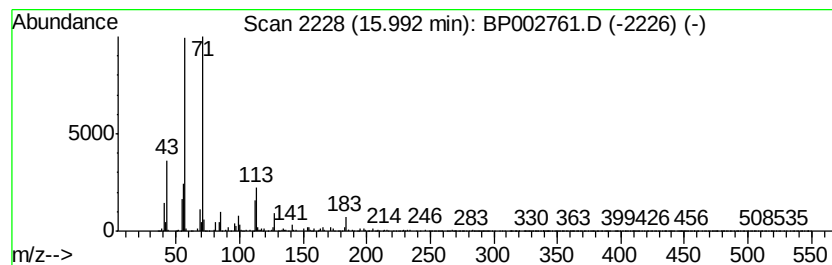
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Heptane, 3-ethyl-5-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.99	2.33 ng	223092	Phenanthrene-d10		16.97	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-5-methyl-		142	C10H22	052896-90-9	72
2	Undecane, 6,6-dimethyl-		184	C13H28	017312-76-4	64
3	Octane, 2,3,7-trimethyl-		156	C11H24	062016-34-6	64
4	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59
5	Tridecane, 7-methyl-		198	C14H30	026730-14-3	59



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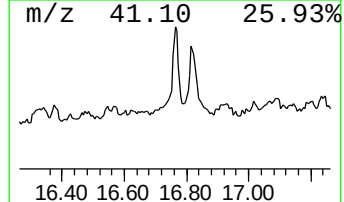
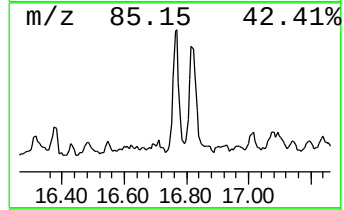
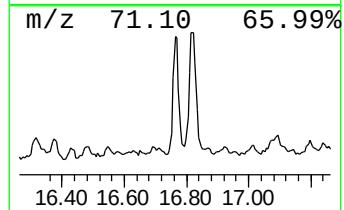
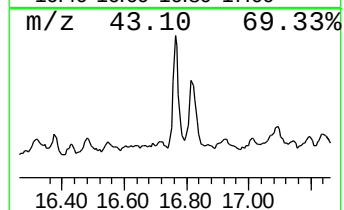
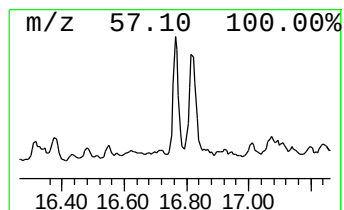
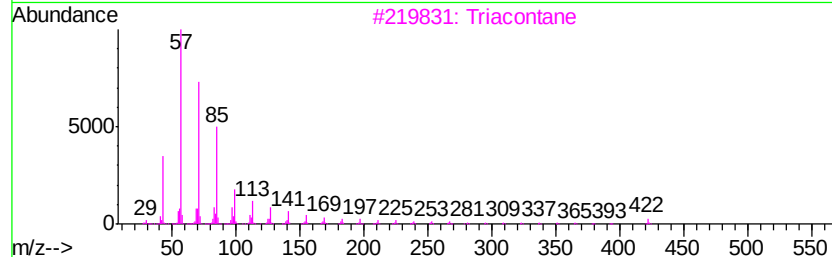
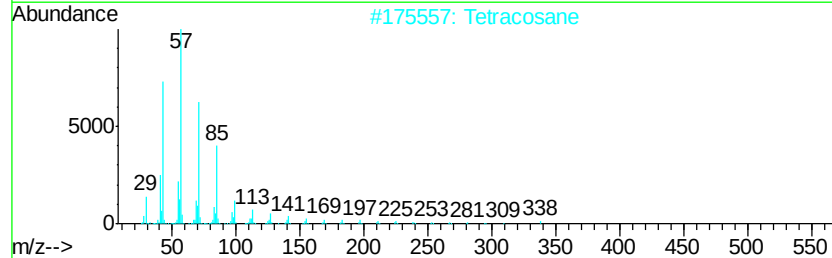
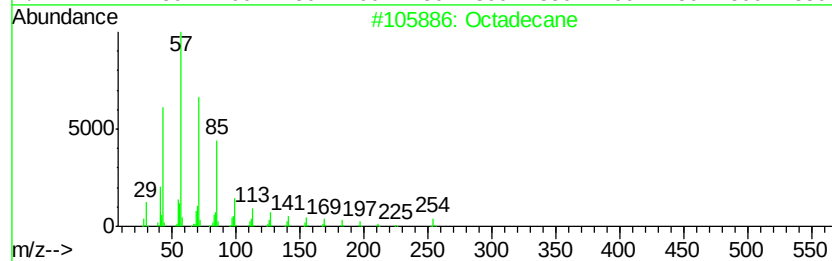
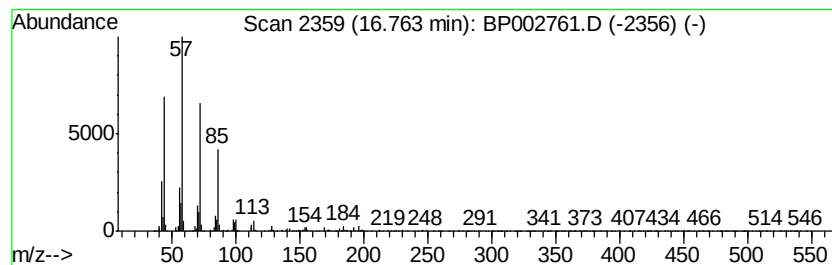
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.76	2.83 ng	270345	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	90
2		Tetracosane	338	C24H50	000646-31-1	87
3		Triacotane	422	C30H62	000638-68-6	87
4		Hexacosane	366	C26H54	000630-01-3	83
5		Octacosane	394	C28H58	000630-02-4	81



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ALS Vial : 13 Sample Multiplier: 1

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FK-GF-02-058-B

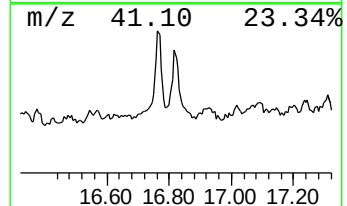
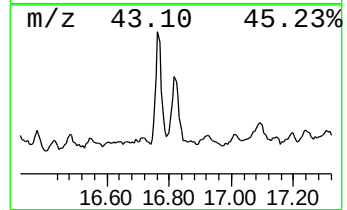
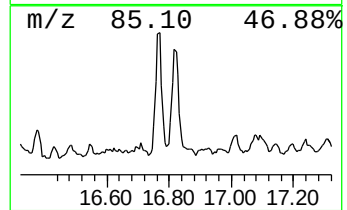
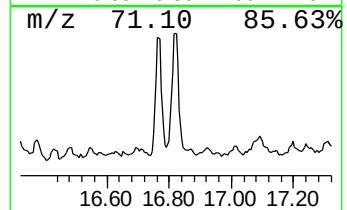
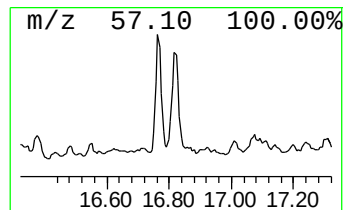
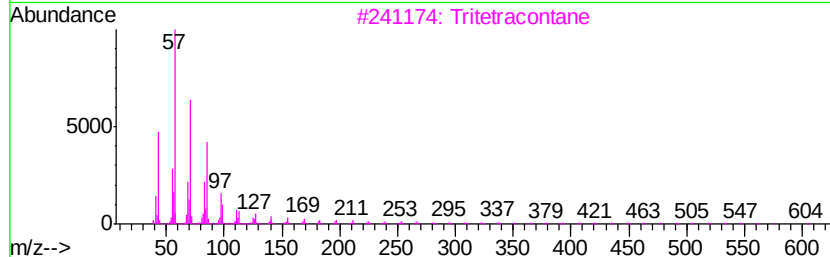
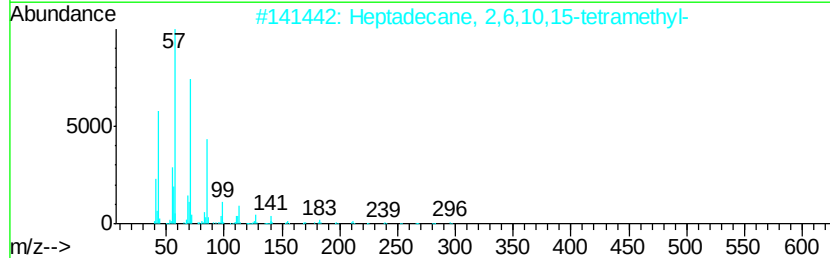
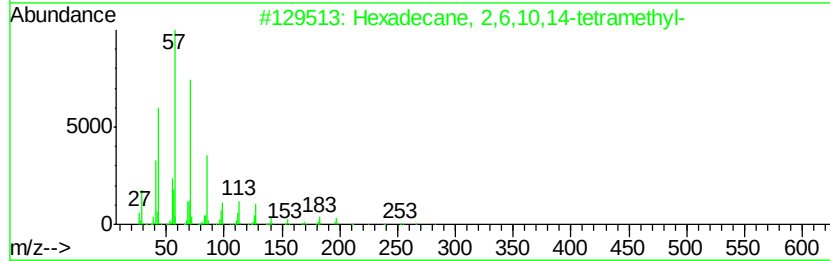
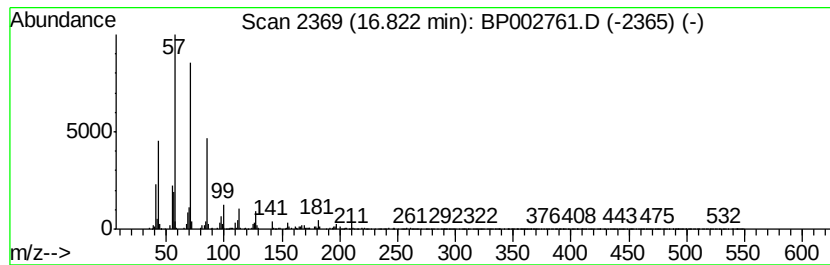
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane, 2,6,10,14-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.82	2.48 ng	236977	Phenanthrene-d10	16.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Tritetracontane	605	C43H88	007098-21-7	90
4		Heptadecane	240	C17H36	000629-78-7	86
5		Nonadecane	268	C19H40	000629-92-5	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
Acq On : 15 Jul 2020 19:03
Operator : CG/JU
Sample : L3304-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FK-GF-02-058-B

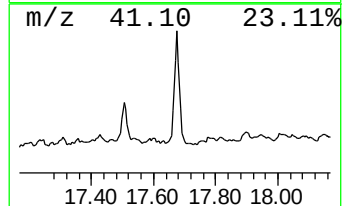
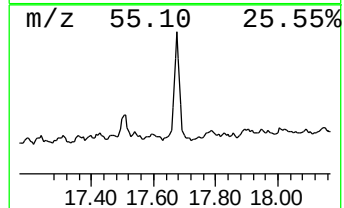
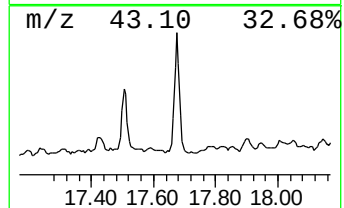
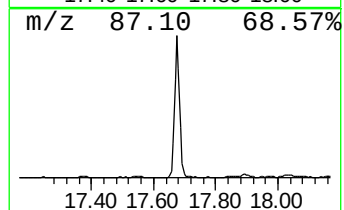
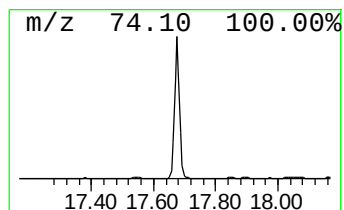
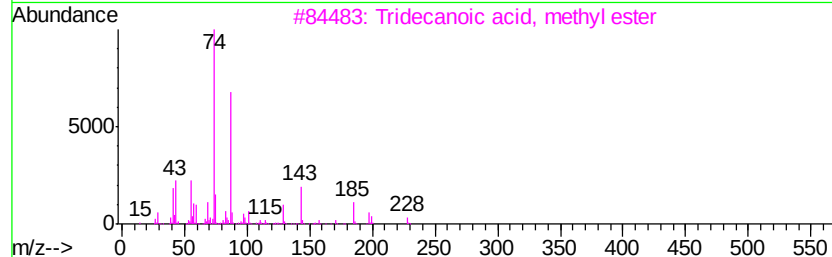
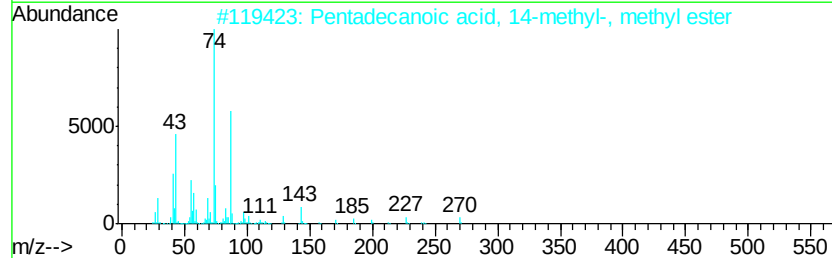
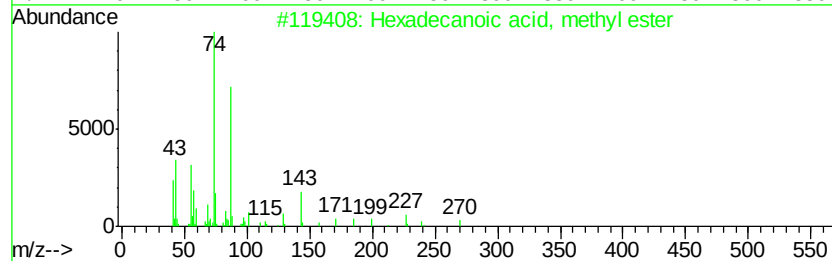
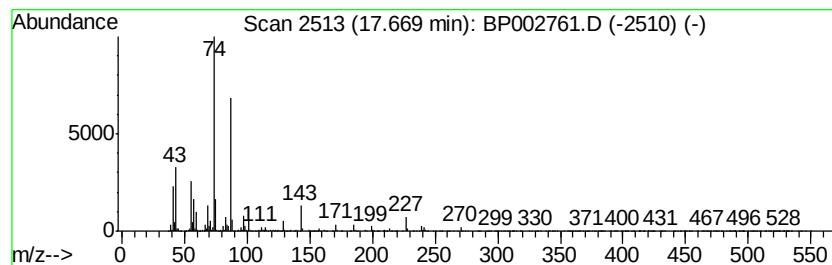
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.67	8.27 ng	791265	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	99
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
Acq On : 15 Jul 2020 19:03
Operator : CG/JU
Sample : L3304-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

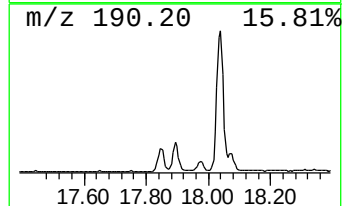
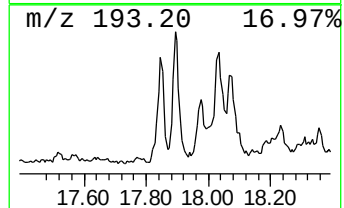
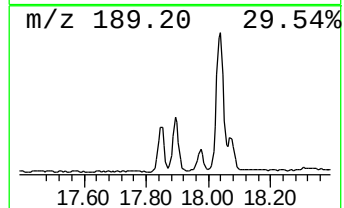
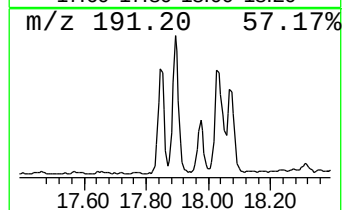
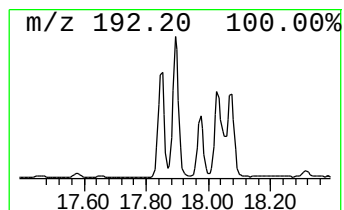
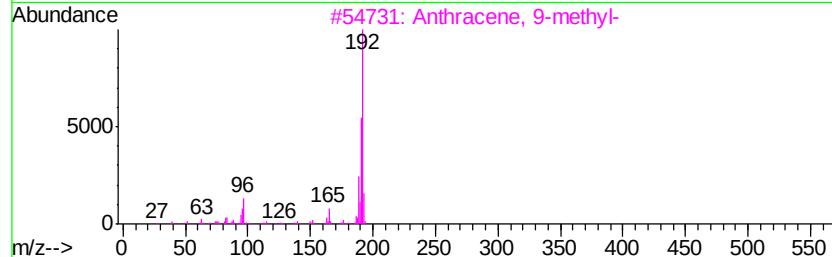
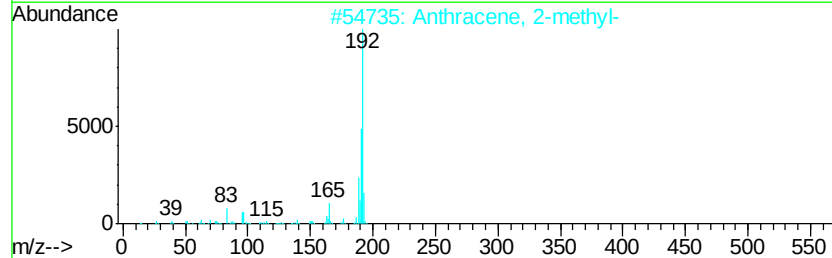
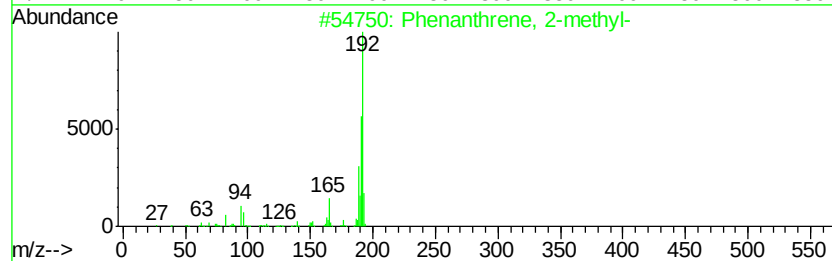
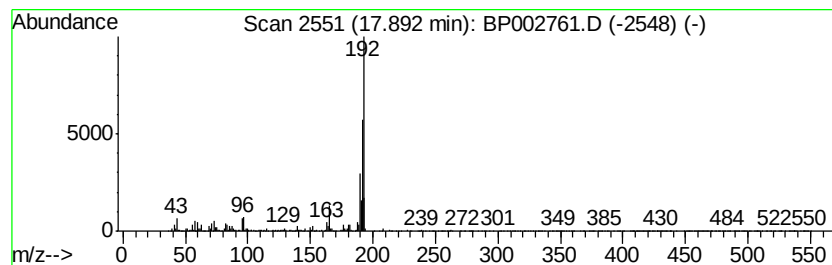
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ClientSampleId :
FK-GF-02-058-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.89	3.35 ng	320888	Phenanthrene-d10		16.97
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
Acq On : 15 Jul 2020 19:03
Operator : CG/JU
Sample : L3304-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

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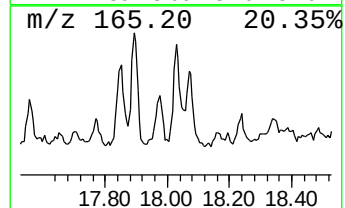
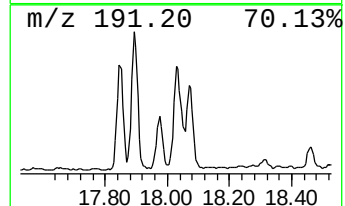
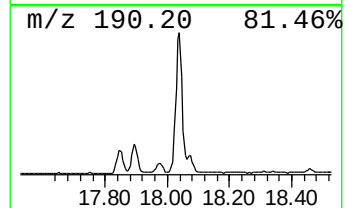
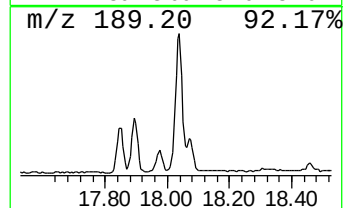
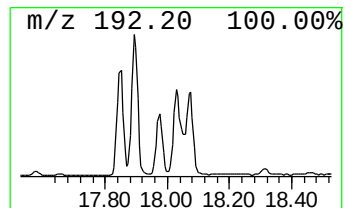
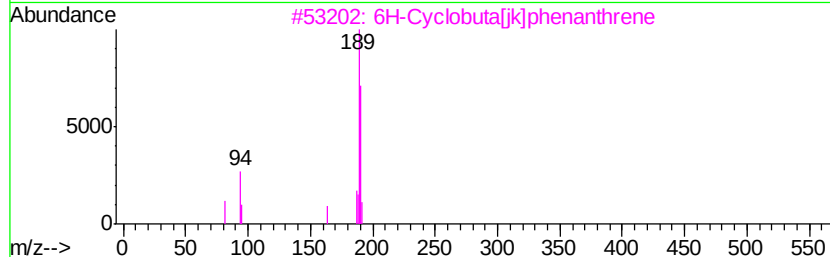
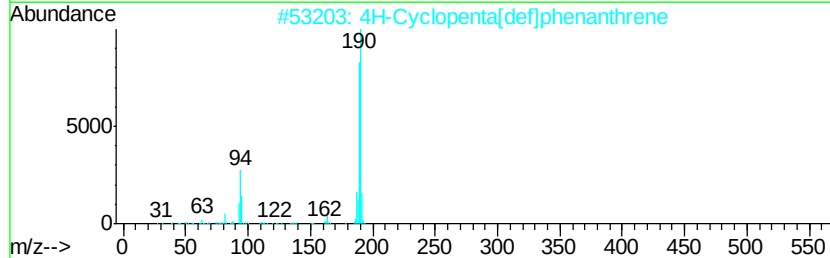
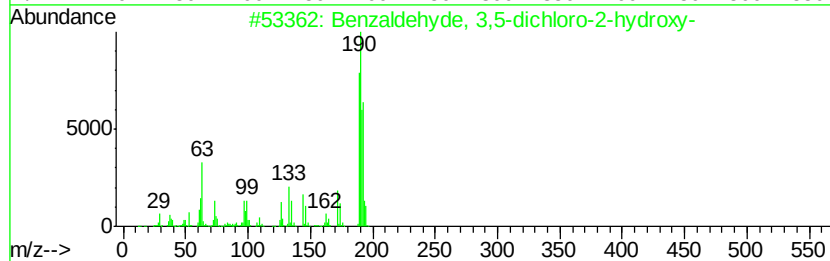
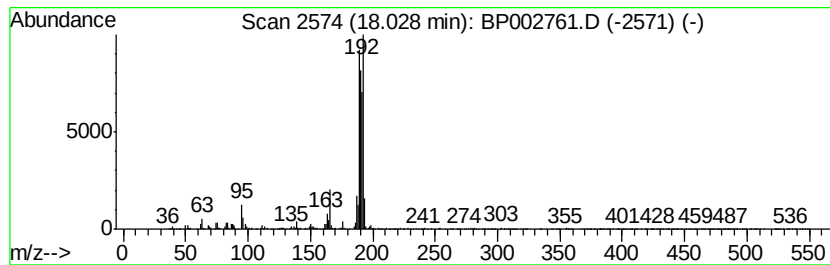
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzaldehyde, 3,5-dichloro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	3.75 ng	358293	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	52
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	30
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
Acq On : 15 Jul 2020 19:03
Operator : CG/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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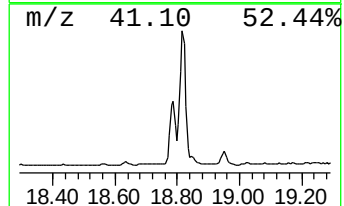
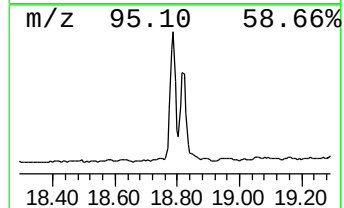
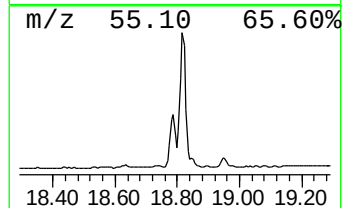
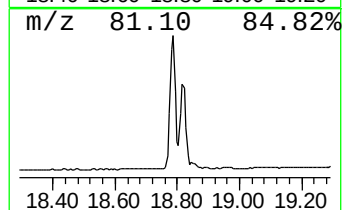
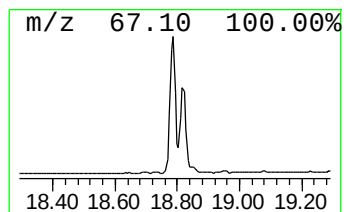
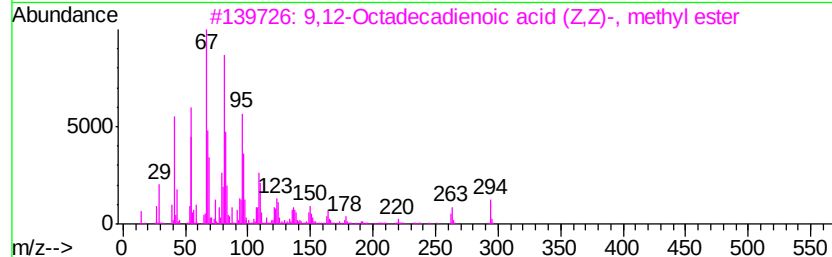
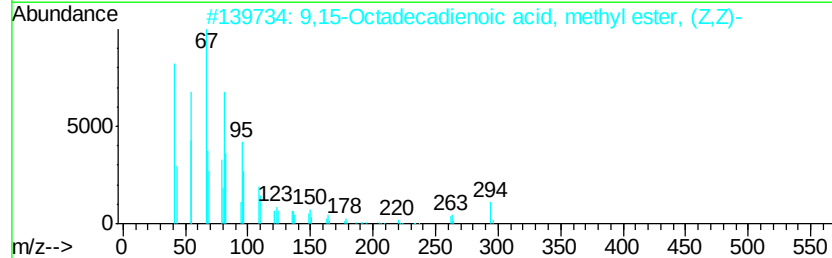
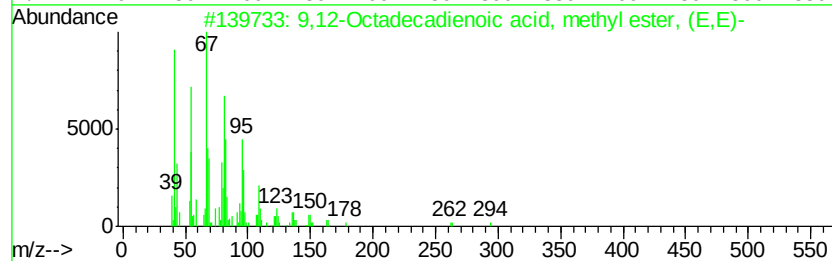
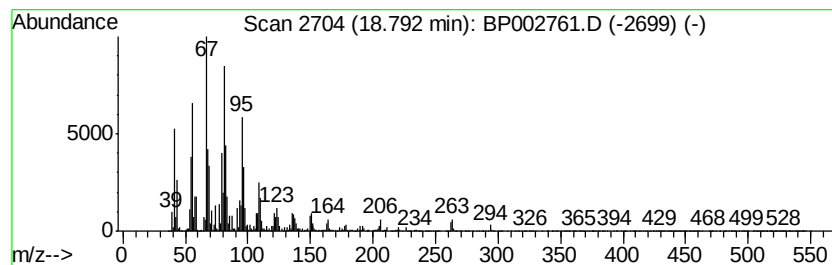
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.79	22.43 ng	2145850	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	98
5		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
Acq On : 15 Jul 2020 19:03
Operator : CG/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FK-GF-02-058-B

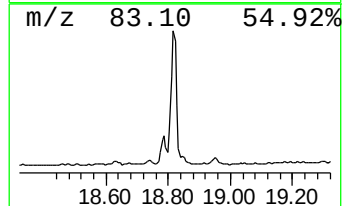
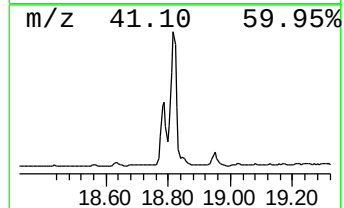
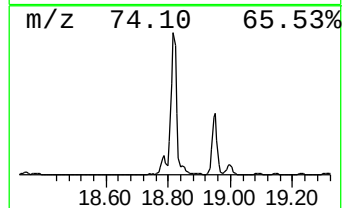
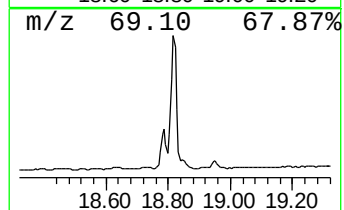
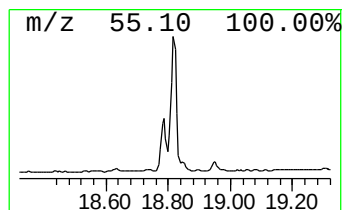
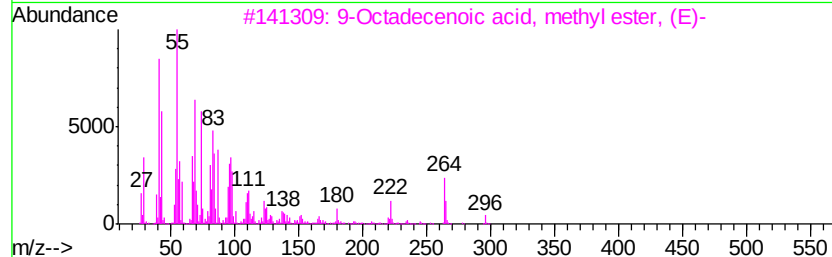
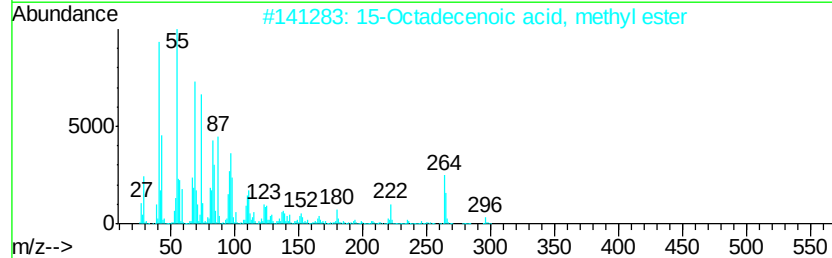
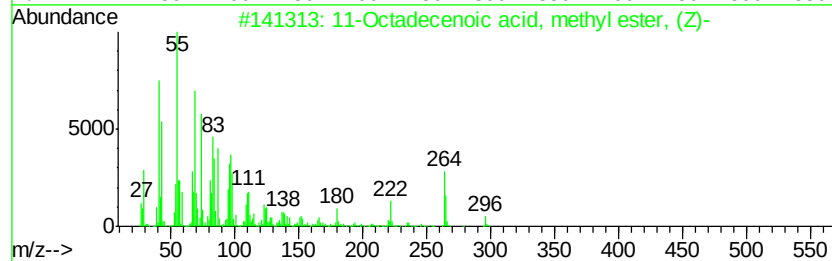
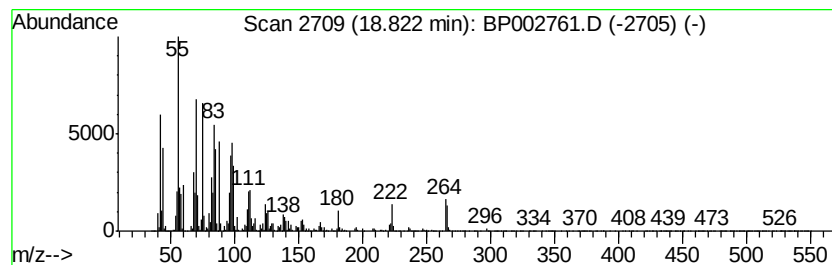
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11-Octadecenoic acid, methyl ester... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	42.36 ng	4051510	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-63-9	99
2		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
3		9-Octadecenoic acid, methyl ester...	296	C19H36O2	001937-62-8	99
4		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
5		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
Data File : BP002761.D
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Sample : L3304-03
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ALS Vial : 13 Sample Multiplier: 1

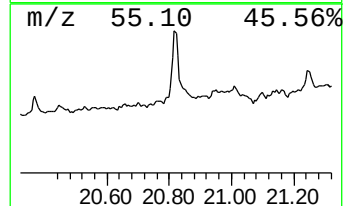
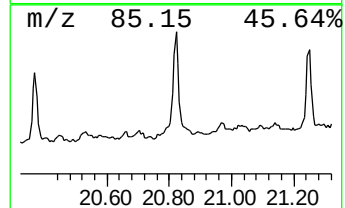
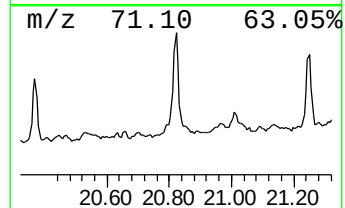
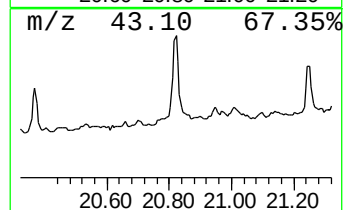
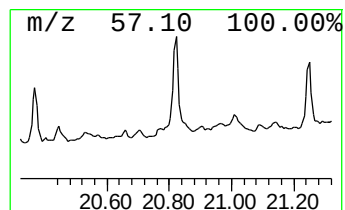
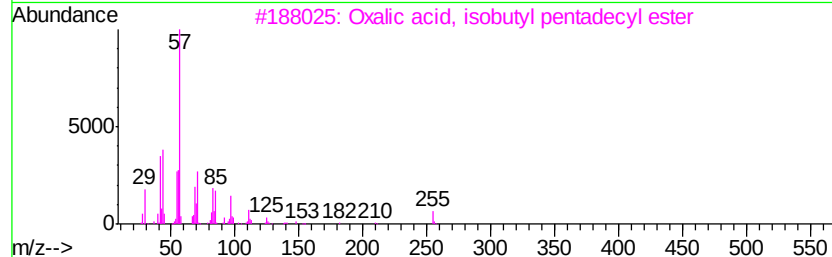
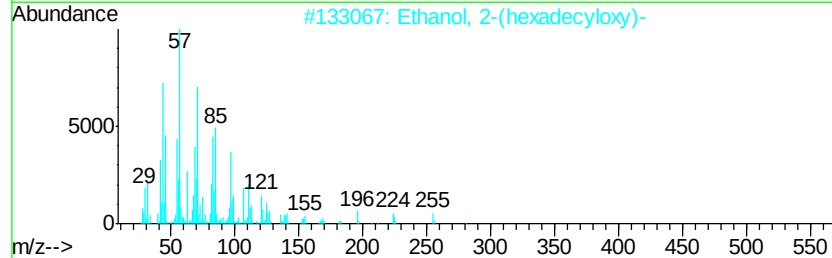
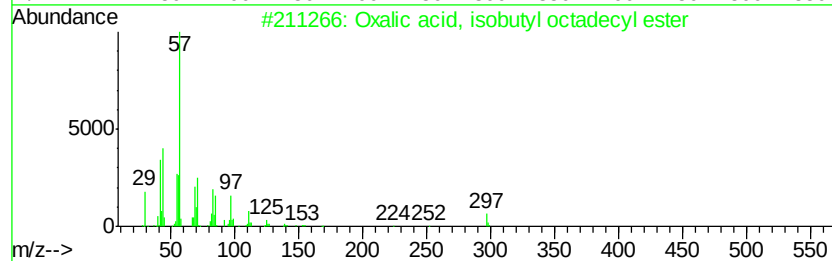
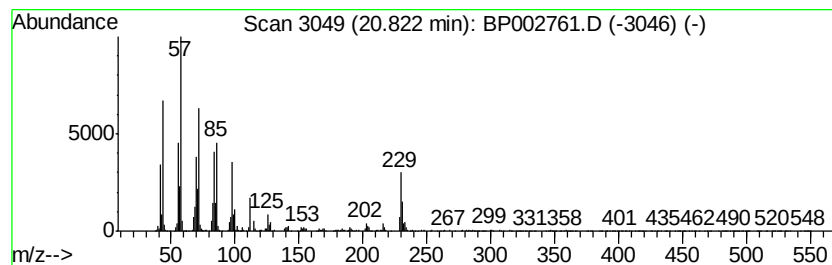
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Oxalic acid, isobutyl octadecyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	4.10 ng	516050	Chrysene-d12	21.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Oxalic acid, isobutyl octadecyl ...	398 C24H46O4	1000309-38-3	74
2	Ethanol, 2-(hexadecyloxy)-	286 C18H38O2	002136-71-2	74
3	Oxalic acid, isobutyl pentadecyl...	356 C21H40O4	1000309-38-0	74
4	Oxalic acid, isobutyl hexadecyl ...	370 C22H42O4	1000309-38-1	74
5	Oxalic acid, isobutyl heptadecyl...	384 C23H44O4	1000309-38-2	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP071520\
 Data File : BP002761.D
 Acq On : 15 Jul 2020 19:03
 Operator : CG/JU
 Sample : L3304-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-GF-02-058-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP071020.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
Hexadecane	15.10	2.1 ng		190267	3	14.21	1813010	20.0
Heptadecane	15.97	2.4 ng		233073	4	16.97	1912980	20.0
Heptane, 3-ethyl-...	15.99	2.3 ng		223092	4	16.97	1912980	20.0
Octadecane	16.76	2.8 ng		270345	4	16.97	1912980	20.0
Hexadecane, 2,6,1...	16.82	2.5 ng		236977	4	16.97	1912980	20.0
Hexadecanoic acid...	17.67	8.3 ng		791265	4	16.97	1912980	20.0
Phenanthrene, 2-m...	17.89	3.4 ng		320888	4	16.97	1912980	20.0
Benzaldehyde, 3,5...	18.03	3.8 ng		358293	4	16.97	1912980	20.0
9,12-Octadecadien...	18.79	22.4 ng		2145850	4	16.97	1912980	20.0
11-Octadecenoic a...	18.82	42.4 ng		4051510	4	16.97	1912980	20.0
Oxalic acid, isob...	20.82	4.1 ng		516050	5	21.08	2517020	20.0