

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
 Data File : BG045796.D
 Acq On : 18 Jun 2020 16:59
 Operator : CG/JU
 Sample : L3029-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WEST-17.5

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_G\METHODS\8270-BG061520.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.533	412	416	438	rVB3	563377	1462623	8.48%	0.422%
2	5.938	478	485	493	rVB2	1148465	2092362	12.13%	0.604%
3	6.561	582	591	595	rBV2	478709	991309	5.75%	0.286%
4	6.919	640	652	657	rBV3	584449	1620006	9.39%	0.468%
5	7.013	664	668	673	rVV	863350	1447112	8.39%	0.418%
6	7.078	673	679	685	rVV3	796056	1838024	10.66%	0.530%
7	7.154	685	692	698	rVV2	903675	1650863	9.57%	0.476%
8	7.313	713	719	724	rBV	557393	983864	5.71%	0.284%
9	7.559	754	761	769	rBV2	2922615	5978948	34.67%	1.726%
10	7.829	802	807	814	rVV3	404437	976832	5.66%	0.282%
11	7.912	815	821	826	rVB2	1047834	1867804	10.83%	0.539%
12	8.217	865	873	877	rBV6	600816	1640797	9.51%	0.474%
13	8.470	910	916	922	rVB2	1308039	2715751	15.75%	0.784%
14	8.570	928	933	948	rVB3	1545903	4125906	23.92%	1.191%
15	8.699	949	955	958	rBV2	970556	1643001	9.53%	0.474%
16	8.887	982	987	991	rBV	589125	1013944	5.88%	0.293%
17	8.934	991	995	1003	rVB	712509	1391786	8.07%	0.402%
18	9.040	1003	1013	1018	rBV2	1554688	3646137	21.14%	1.052%
19	9.175	1030	1036	1044	rVB	3928291	6880298	39.90%	1.986%
20	9.292	1044	1056	1062	rBV7	1155455	3452040	20.02%	0.996%
21	9.374	1062	1070	1072	rBV3	576482	1224379	7.10%	0.353%
22	9.574	1099	1104	1109	rBV2	485910	977334	5.67%	0.282%
23	9.633	1109	1114	1115	rVV	730860	1113219	6.46%	0.321%
24	9.656	1115	1118	1124	rVB2	1375521	2227427	12.92%	0.643%
25	9.821	1137	1146	1150	rBV2	712658	1618034	9.38%	0.467%
26	9.868	1150	1154	1158	rVB3	692729	1063185	6.17%	0.307%
27	9.921	1158	1163	1170	rBV5	919839	2025420	11.74%	0.585%
28	9.985	1170	1174	1176	rBV	658772	997304	5.78%	0.288%
29	10.021	1177	1180	1185	rVB2	1281026	1939451	11.25%	0.560%
30	10.091	1185	1192	1195	rBV	1603932	2792412	16.19%	0.806%
31	10.138	1195	1200	1203	rBV2	859402	1438906	8.34%	0.415%
32	10.273	1218	1223	1227	rBV2	851657	1508625	8.75%	0.435%
33	10.367	1235	1239	1245	rVB	1420750	2522296	14.63%	0.728%
34	10.438	1246	1251	1256	rBV	783192	1348762	7.82%	0.389%

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 Sampling : 1
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 Stop Thrs : 0
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 Min Area: 3 % of largest Peak
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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061520.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.526	1260	1266	1271	rVB3	711758	1328701	7.70%	0.383%
36	10.649	1282	1287	1291	rBV4	653791	1170355	6.79%	0.338%
37	10.731	1291	1301	1307	rVB2	5629662	11475912	66.55%	3.312%
38	10.790	1307	1311	1314	rBV4	520018	925217	5.37%	0.267%
39	10.861	1319	1323	1327	rVV3	1164253	2393686	13.88%	0.691%
40	10.914	1327	1332	1336	rVV	1731789	2861710	16.59%	0.826%
41	10.966	1336	1341	1349	rVV6	817978	1790131	10.38%	0.517%
42	11.207	1378	1382	1391	rVB2	1261431	2834435	16.44%	0.818%
43	11.319	1396	1401	1405	rBV3	666161	1093989	6.34%	0.316%
44	11.407	1411	1416	1418	rBV2	735190	1251589	7.26%	0.361%
45	11.448	1418	1423	1431	rVV5	1474052	3646721	21.15%	1.053%
46	11.513	1431	1434	1440	rVB3	928464	1585094	9.19%	0.457%
47	11.589	1441	1447	1453	rBV3	1836978	3194865	18.53%	0.922%
48	11.665	1453	1460	1466	rBV3	2064916	4384864	25.43%	1.266%
49	11.777	1473	1479	1488	rBV3	2315517	5086505	29.50%	1.468%
50	11.853	1489	1492	1498	rVB3	787253	1401126	8.12%	0.404%
51	11.942	1499	1507	1514	rBV2	3153875	6297422	36.52%	1.818%
52	12.182	1541	1548	1553	rBV	6047748	10516532	60.98%	3.035%
53	12.294	1562	1567	1572	rVB2	1404787	2740790	15.89%	0.791%
54	12.394	1577	1584	1592	rVB6	1528332	4117891	23.88%	1.188%
55	12.617	1617	1622	1624	rBV2	737148	1236020	7.17%	0.357%
56	12.664	1624	1630	1634	rVV2	2811265	5228155	30.32%	1.509%
57	12.799	1650	1653	1657	rVB4	1088378	1603920	9.30%	0.463%
58	12.846	1657	1661	1668	rVB3	1159891	2298619	13.33%	0.663%
59	12.911	1668	1672	1676	rBV3	1473279	2585510	14.99%	0.746%
60	12.981	1679	1684	1690	rVB3	1263078	2608728	15.13%	0.753%
61	13.069	1694	1699	1705	rBV3	1647441	3331867	19.32%	0.962%
62	13.128	1705	1709	1714	rVB2	3360803	4483991	26.00%	1.294%
63	13.287	1732	1736	1742	rVB5	933481	1446197	8.39%	0.417%
64	13.428	1752	1760	1766	rVV	7693511	13805479	80.05%	3.985%
65	13.534	1771	1778	1782	rVV2	2181641	5058733	29.33%	1.460%
66	13.581	1782	1786	1790	rVB3	1426536	1961765	11.38%	0.566%
67	13.745	1809	1814	1821	rBV6	937760	2006216	11.63%	0.579%
68	13.863	1830	1834	1840	rBV5	1080745	2985364	17.31%	0.862%
69	14.050	1862	1866	1867	rBV3	1670310	2075784	12.04%	0.599%
70	14.080	1868	1871	1874	rVV2	2628388	3817171	22.13%	1.102%
71	14.115	1874	1877	1882	rVB3	2018189	2688675	15.59%	0.776%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	14.509	1936	1944	1947	rBV	9183837	15216377	88.24%	4.392%
73	14.562	1949	1953	1966	rVV7	1565479	4154192	24.09%	1.199%
74	14.955	2011	2020	2023	rBV6	1535382	4172779	24.20%	1.204%
75	14.990	2023	2026	2032	rVB2	1778566	2619908	15.19%	0.756%
76	15.049	2033	2036	2041	rVV2	826131	1461373	8.47%	0.422%
77	15.114	2043	2047	2055	rVB4	2624560	4474233	25.94%	1.291%
78	15.178	2055	2058	2060	rBV	827883	942141	5.46%	0.272%
79	15.378	2089	2092	2096	rBV2	782572	944500	5.48%	0.273%
80	15.460	2099	2106	2109	rVB	9146329	14047629	81.46%	4.054%
81	15.678	2139	2143	2148	rBV6	814920	1470278	8.53%	0.424%
82	15.854	2168	2173	2177	rVB	4126751	6778768	39.31%	1.956%
83	16.001	2194	2198	2203	rBV3	1392302	2497497	14.48%	0.721%
84	16.060	2204	2208	2218	rVB5	1911170	3816202	22.13%	1.101%
85	16.318	2245	2252	2255	rBV	9457058	17245222	100.00%	4.977%
86	16.764	2325	2328	2332	rBV2	753226	993289	5.76%	0.287%
87	16.911	2350	2353	2356	rVB2	893870	1087127	6.30%	0.314%
88	17.111	2380	2387	2390	rVV	8915482	13442490	77.95%	3.880%
89	17.158	2390	2395	2404	rVB	4521733	7770208	45.06%	2.243%
90	17.340	2422	2426	2429	rBV	898745	1320006	7.65%	0.381%
91	17.634	2473	2476	2482	rVB3	1060579	1519327	8.81%	0.439%
92	17.751	2492	2496	2505	rVV4	1094526	2129903	12.35%	0.615%
93	17.845	2506	2512	2516	rVB	8049542	11032494	63.97%	3.184%
94	18.116	2554	2558	2566	rBV3	730469	1448481	8.40%	0.418%
95	18.527	2623	2628	2633	rBV	6478049	8612458	49.94%	2.486%
96	19.155	2730	2735	2739	rVB	5061548	6423587	37.25%	1.854%
97	19.725	2828	2832	2836	rVB	3624225	3978091	23.07%	1.148%
98	19.942	2865	2869	2874	rVB	1825462	2451585	14.22%	0.708%
99	20.254	2917	2922	2926	rVB	2499524	2937445	17.03%	0.848%
100	20.741	3001	3005	3016	rVB	1536216	1954693	11.33%	0.564%

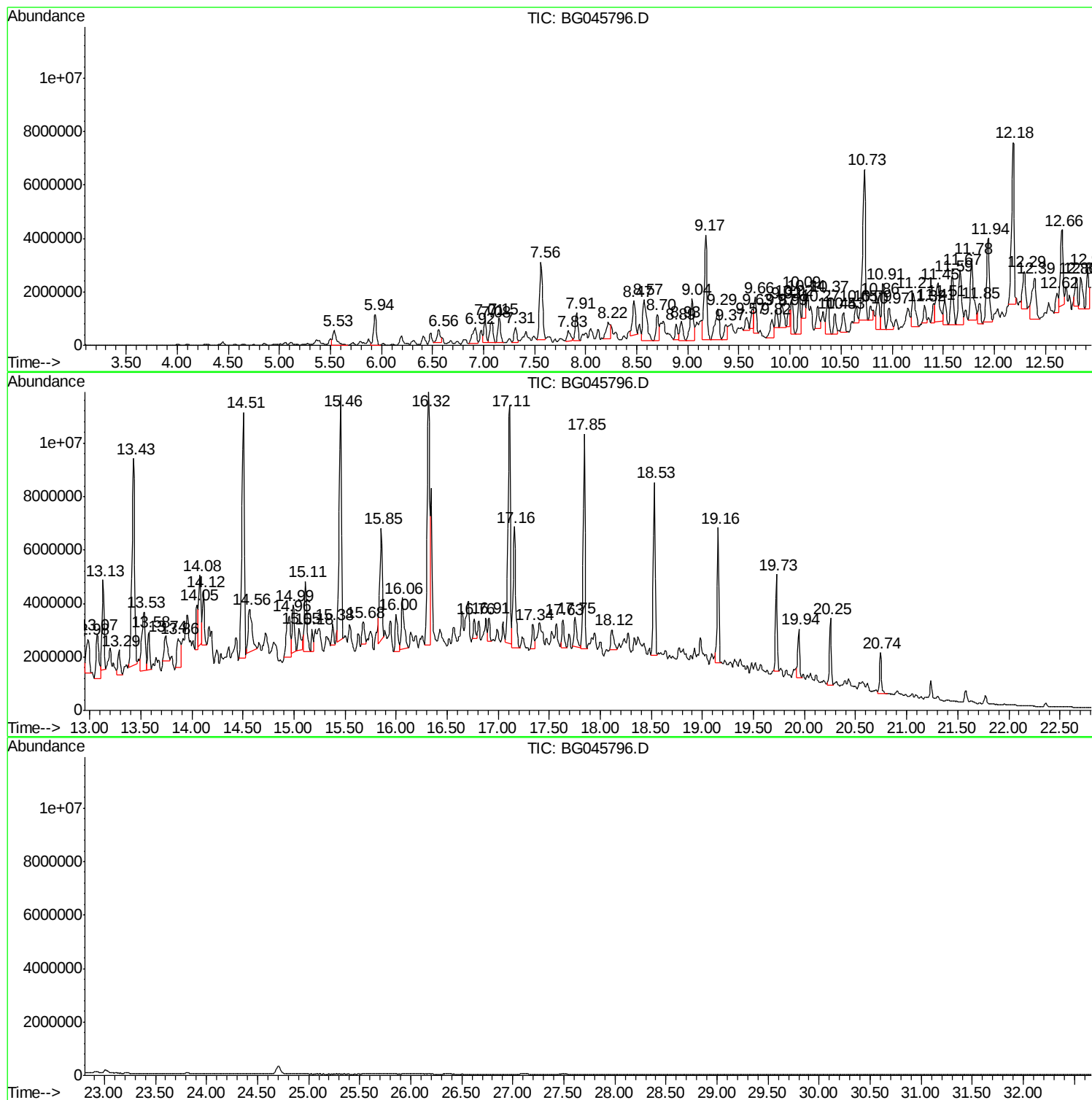
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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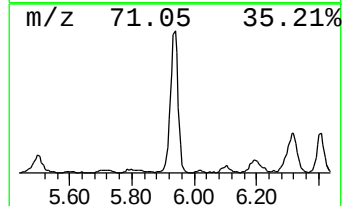
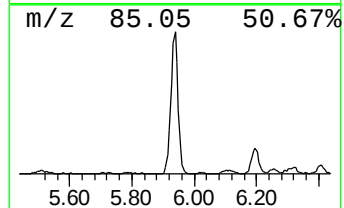
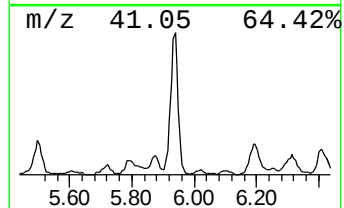
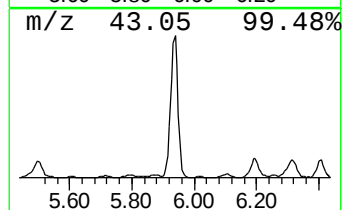
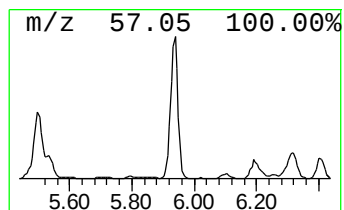
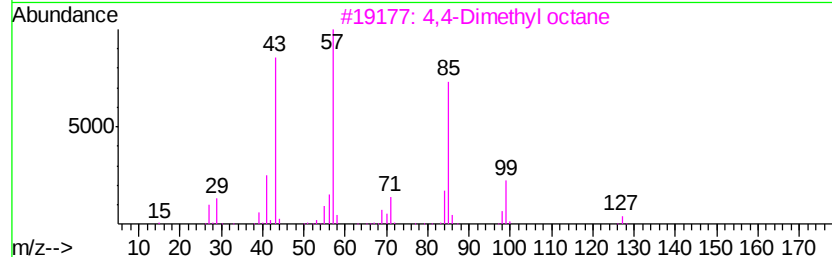
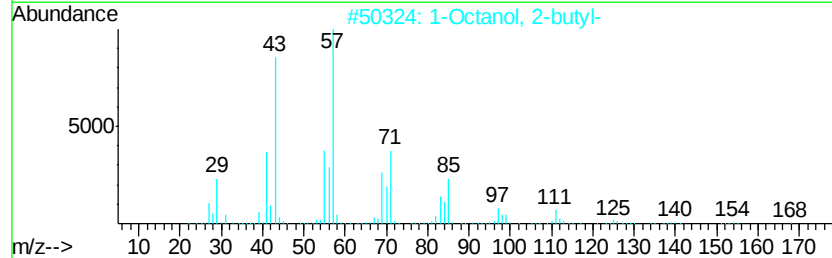
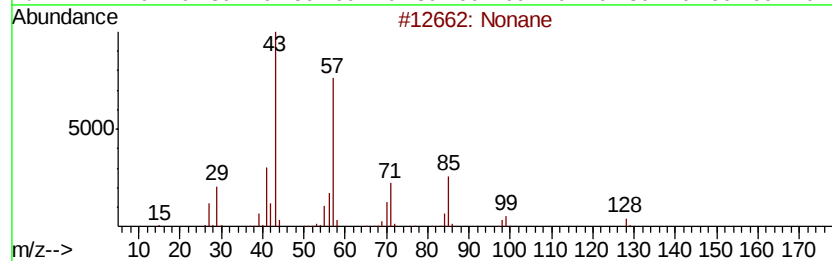
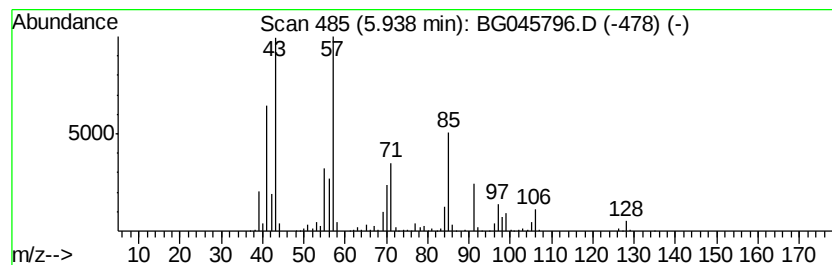
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Nonane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.94	22.40 ng	2092360	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane	128	C9H20	000111-84-2	70
2		1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	64
3		4,4-Dimethyl octane	142	C10H22	015869-95-1	47
4		Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	43
5		Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1000309-37-4	38



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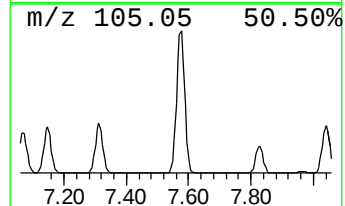
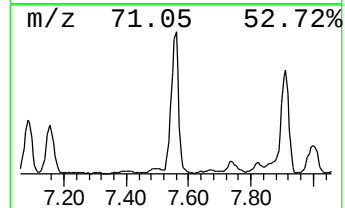
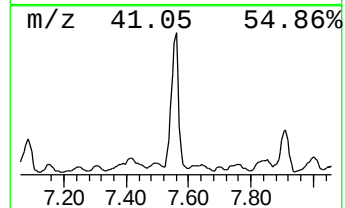
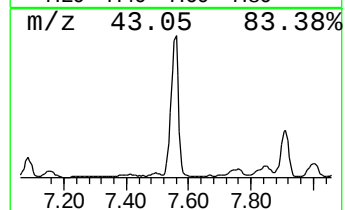
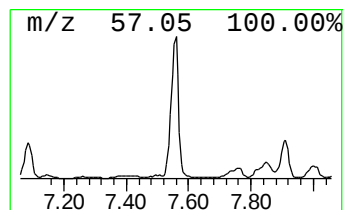
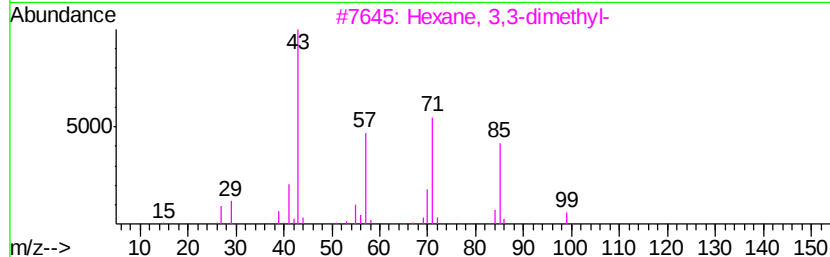
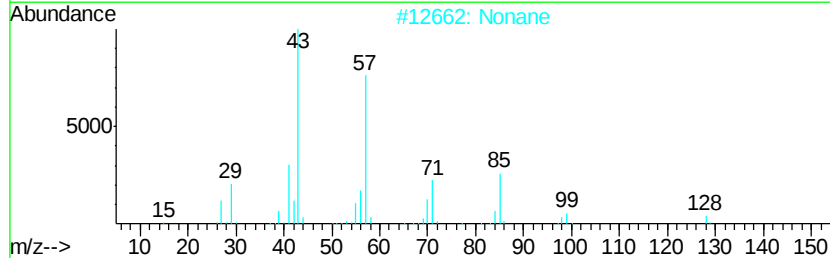
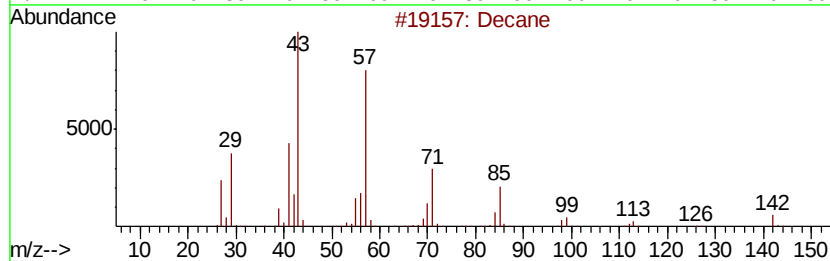
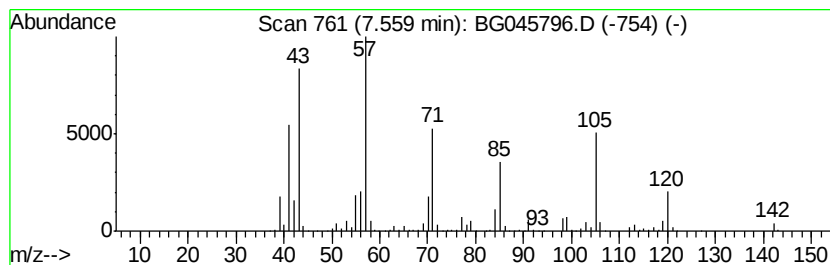
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Decane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	64.02 ng	5978950	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	93
2		Nonane	128	C9H20	000111-84-2	46
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	43
4		Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	43
5		Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	38



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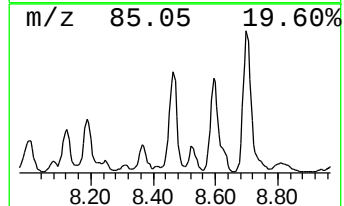
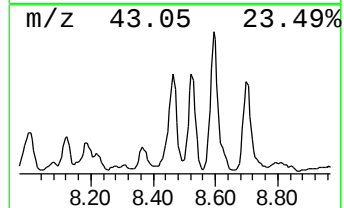
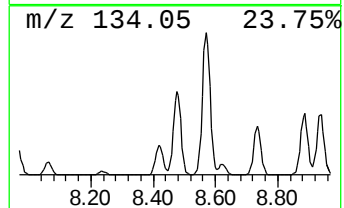
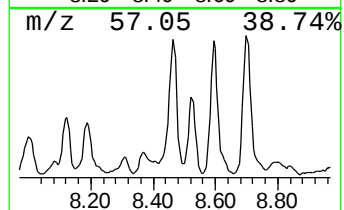
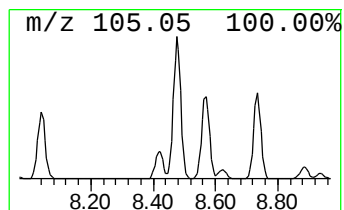
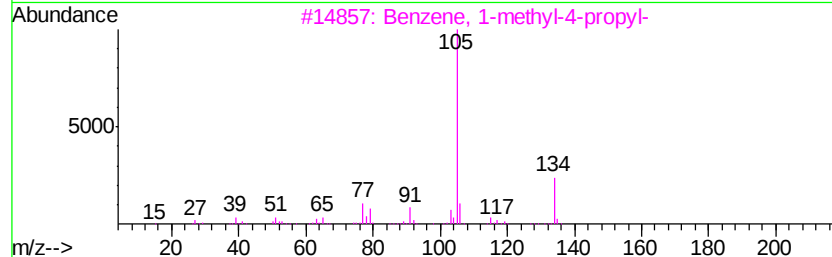
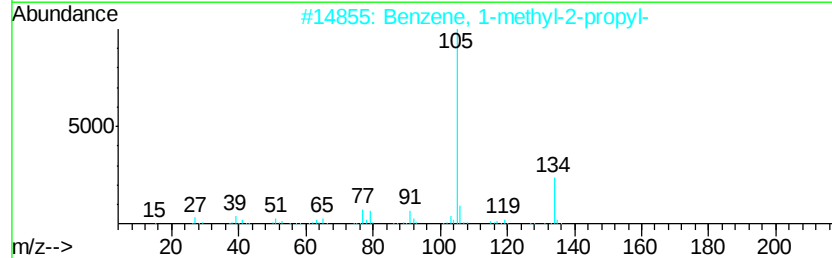
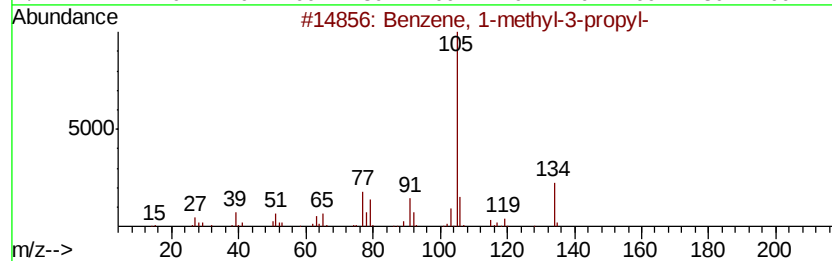
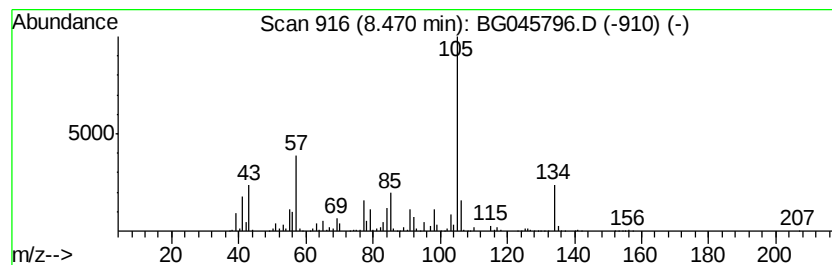
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-methyl-3-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.47	29.08 ng	2715750	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	60
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	49
3		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	49
4		Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	46
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
Sample : L3029-01
Misc :
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ClientSampleId :
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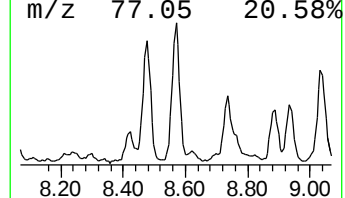
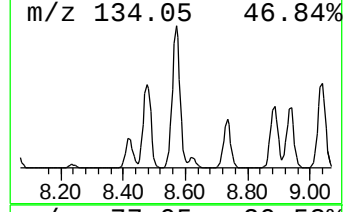
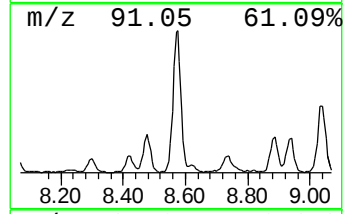
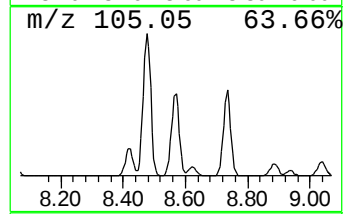
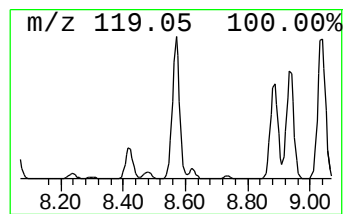
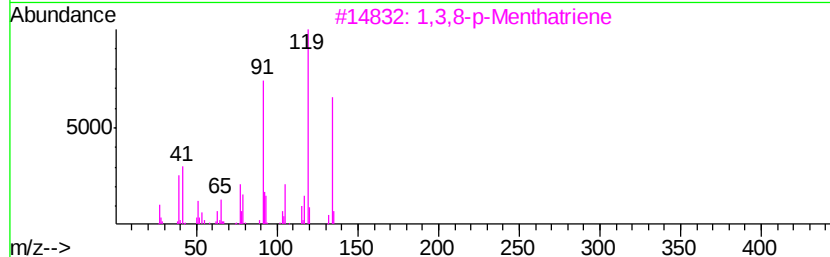
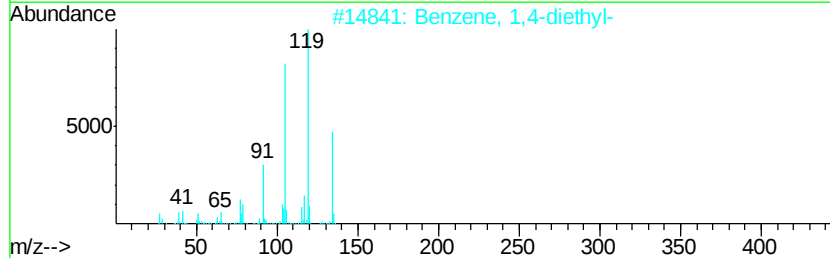
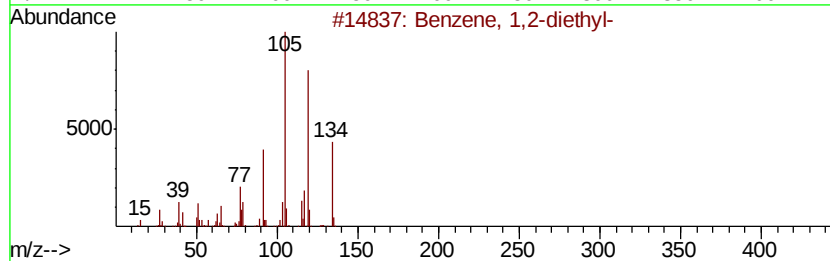
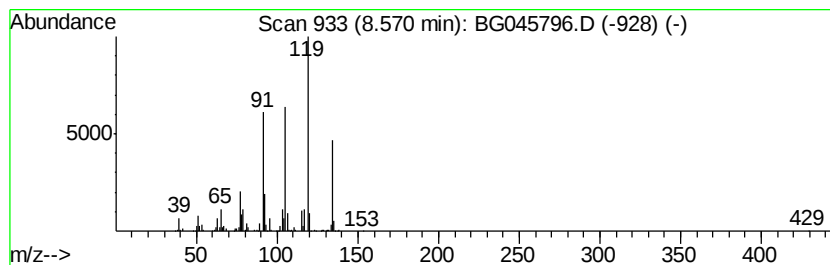
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, 1,2-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.57	44.18 ng	4125910	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	93
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
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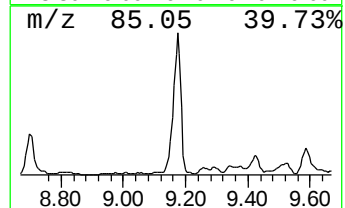
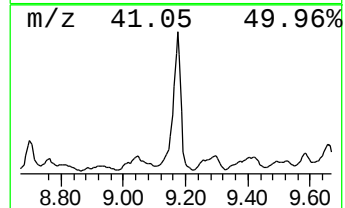
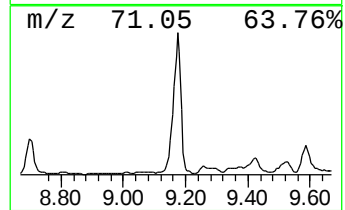
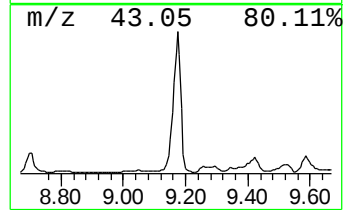
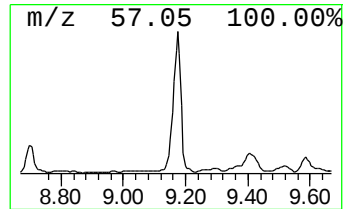
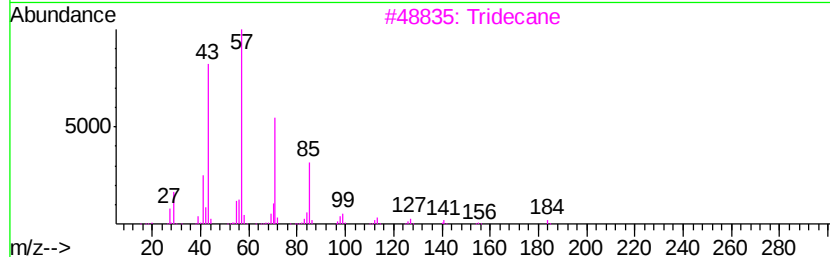
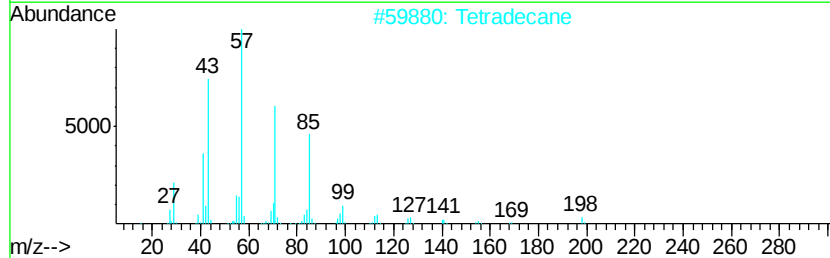
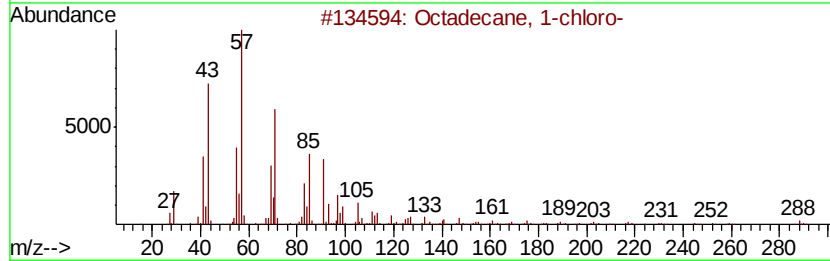
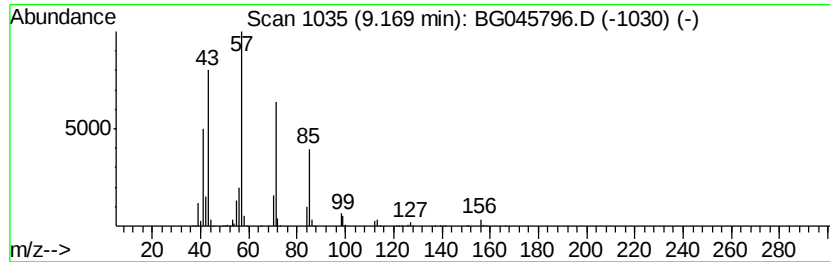
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octadecane, 1-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.17	73.67 ng	6880300	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	78
2		Tetradecane	198	C14H30	000629-59-4	72
3		Tridecane	184	C13H28	000629-50-5	72
4		Undecane	156	C11H24	001120-21-4	70
5		Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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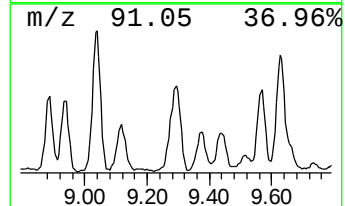
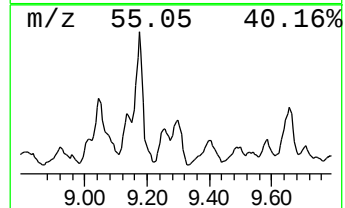
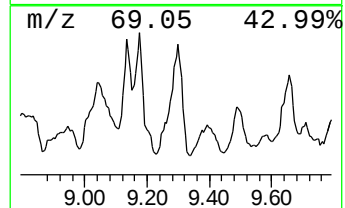
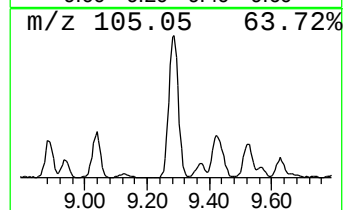
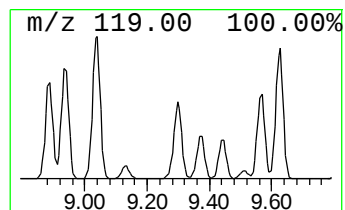
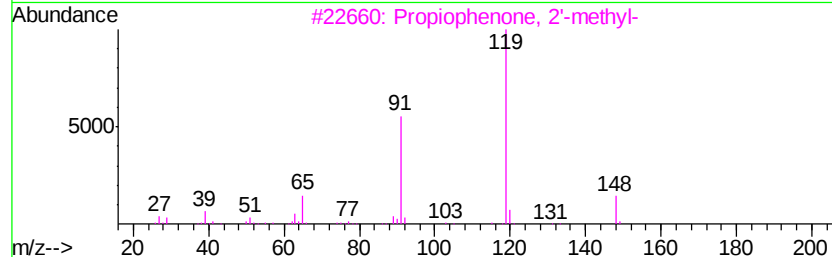
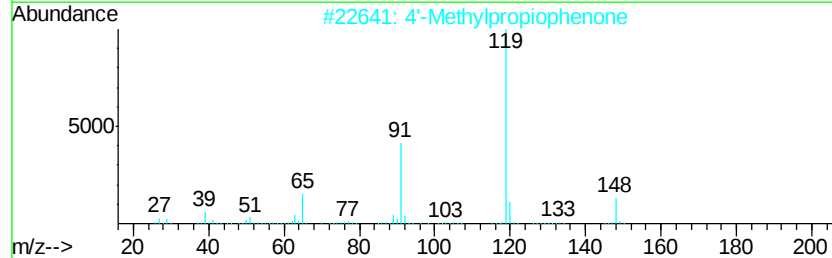
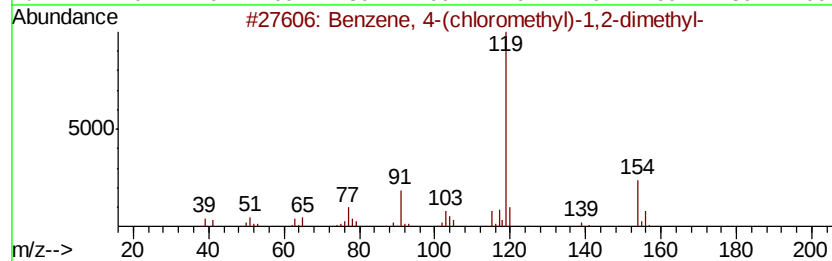
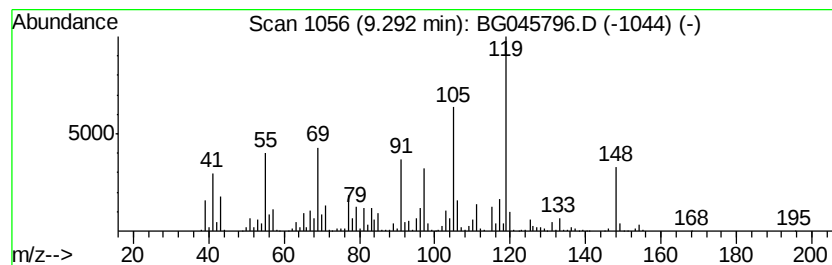
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 4-(chloromethyl)-1... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.29	36.96 ng	3452040	1,4-Dichlorobenzene-d4	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-(chloromethyl)-1,2-di...	154	C9H11Cl	000102-46-5	55
2		4'-Methylpropiophenone	148	C10H12O	005337-93-9	49
3		Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	49
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	49
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
Sample : L3029-01
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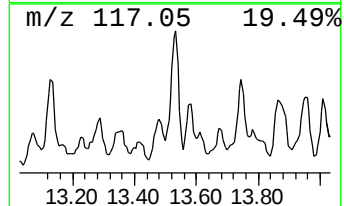
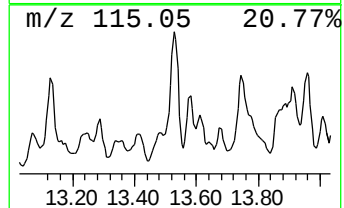
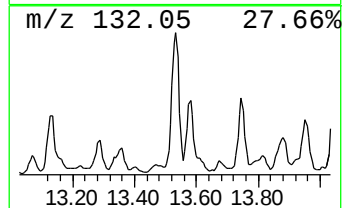
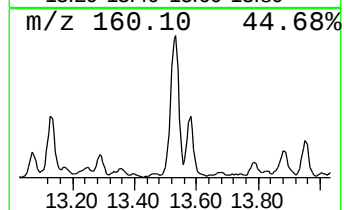
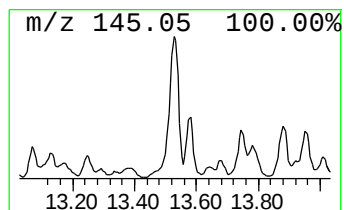
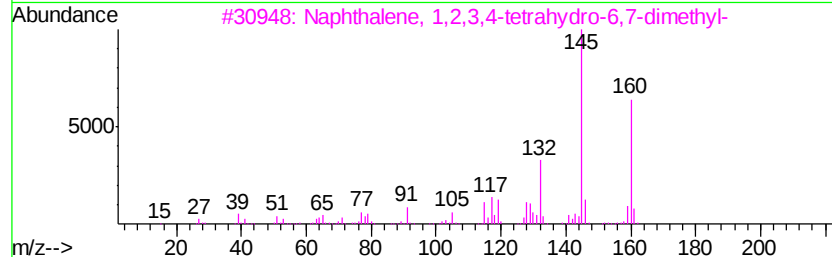
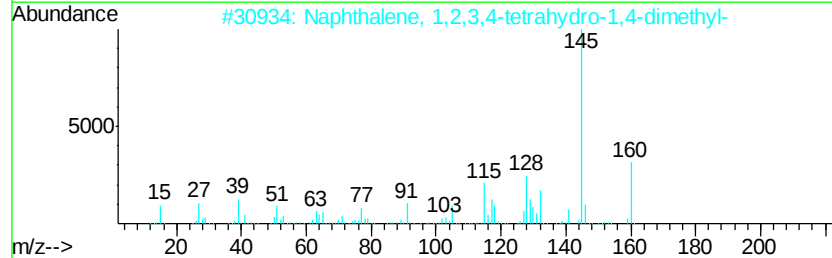
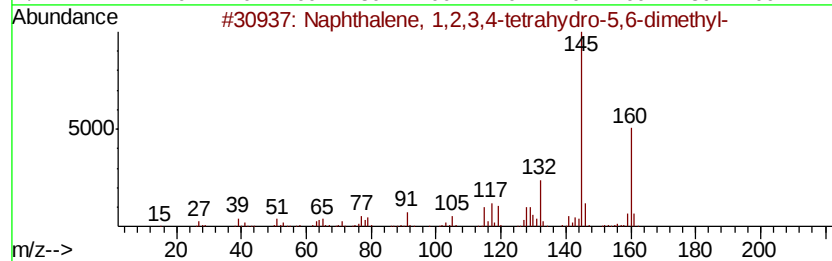
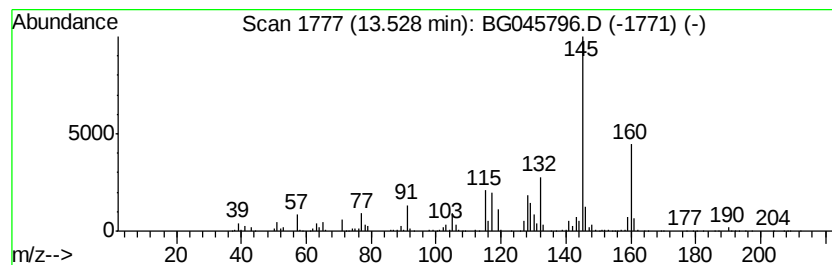
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	24.35 ng	5058730	Acenaphthene-d10	14.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	020027-77-4	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	93
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	93
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	91
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	90



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Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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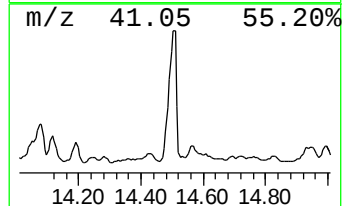
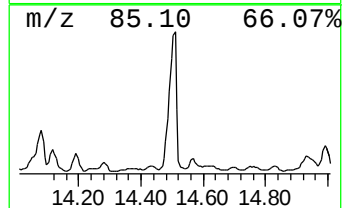
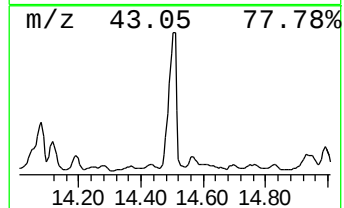
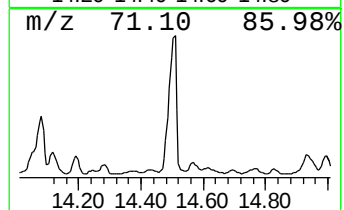
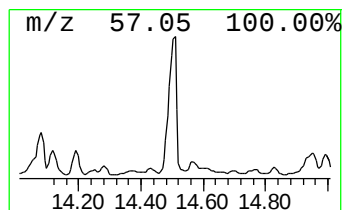
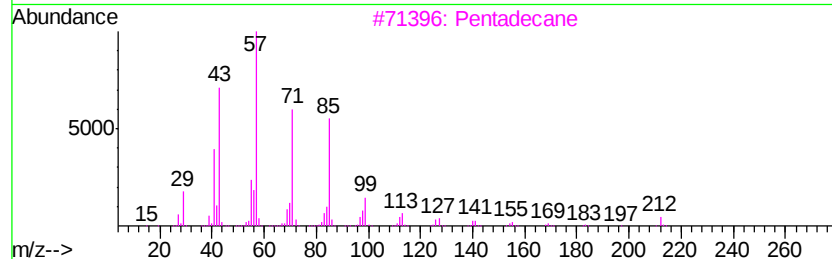
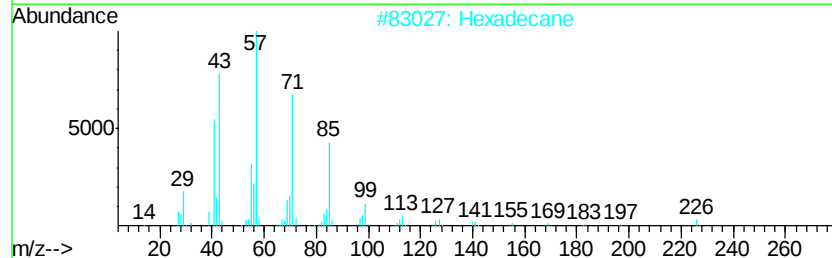
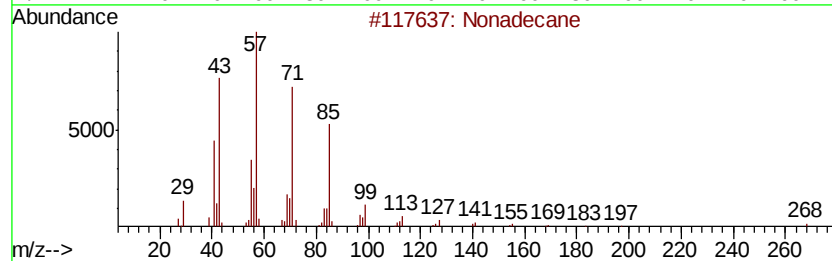
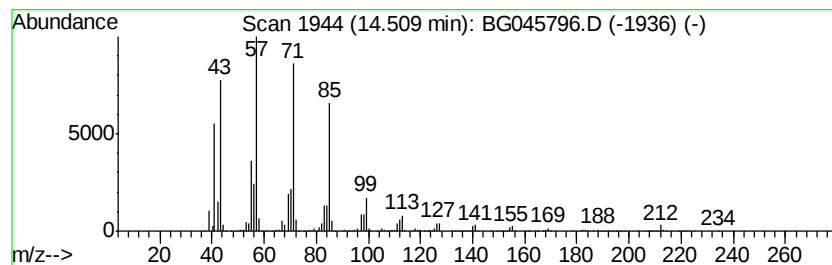
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Nonadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.51	73.26 ng	15216400	Acenaphthene-d10	14.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Hexadecane	226	C16H34	000544-76-3	91
3		Pentadecane	212	C15H32	000629-62-9	91
4		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86
5		Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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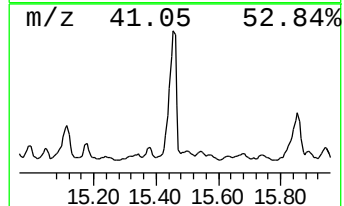
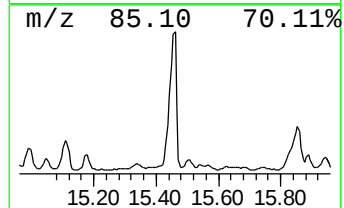
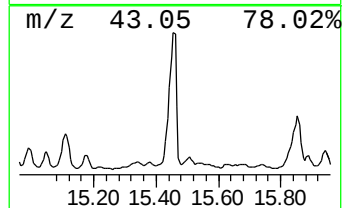
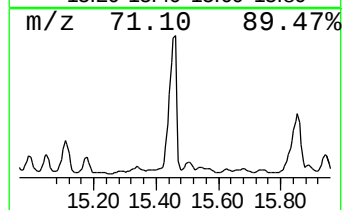
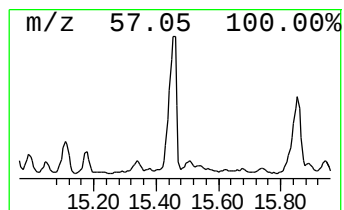
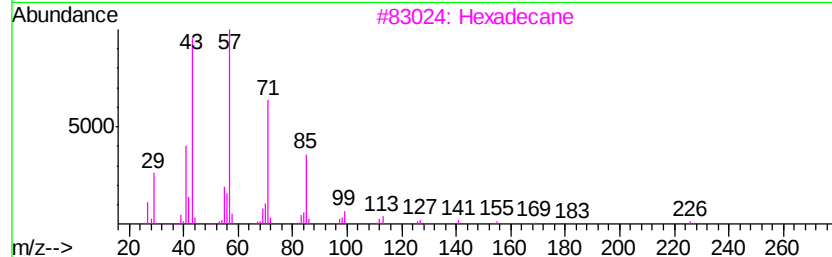
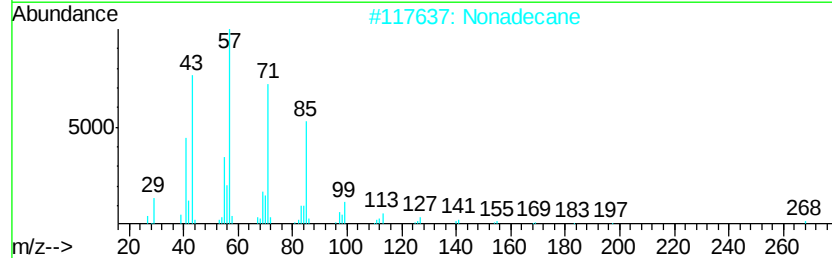
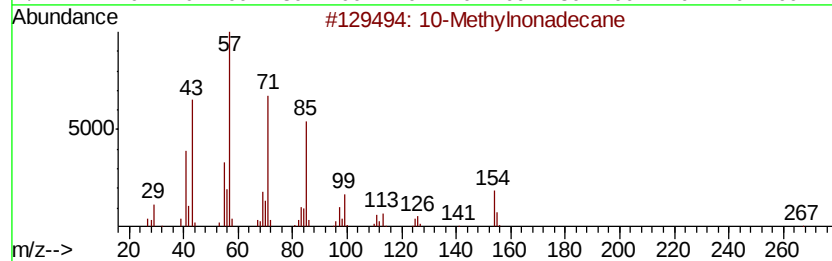
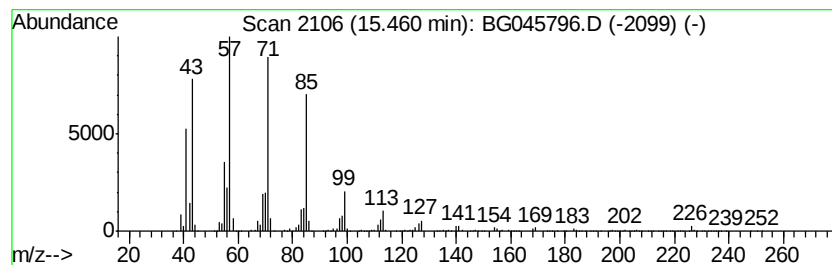
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 10-Methylnonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	67.63 ng	14047600	Acenaphthene-d10	14.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Methylnonadecane	282	C20H42	056862-62-5	91
2		Nonadecane	268	C19H40	000629-92-5	90
3		Hexadecane	226	C16H34	000544-76-3	87
4		Heptadecane	240	C17H36	000629-78-7	86
5		Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
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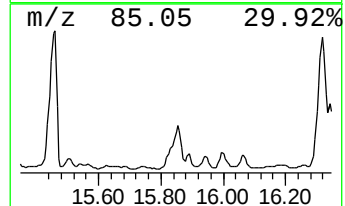
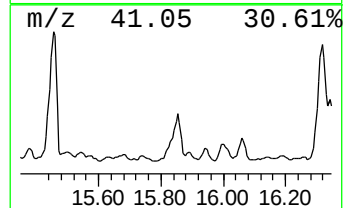
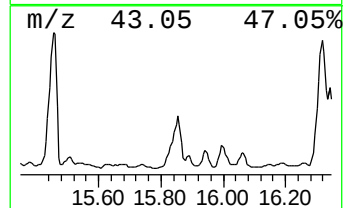
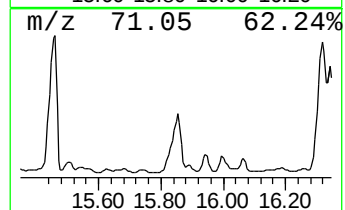
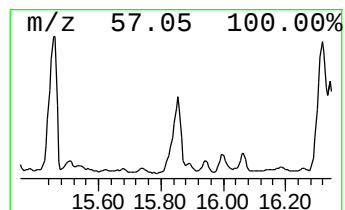
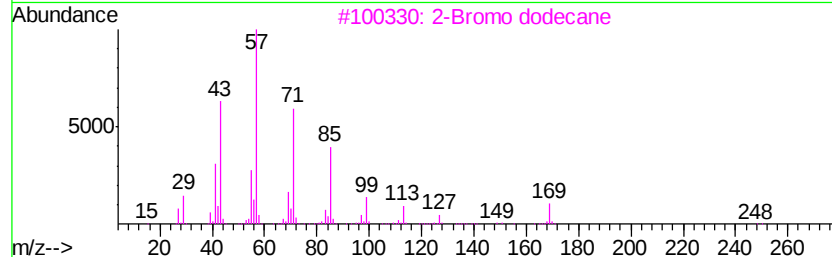
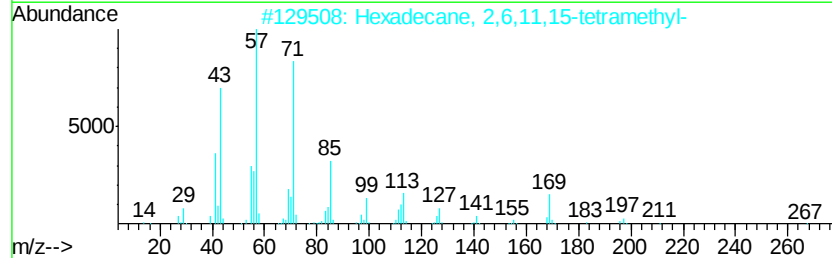
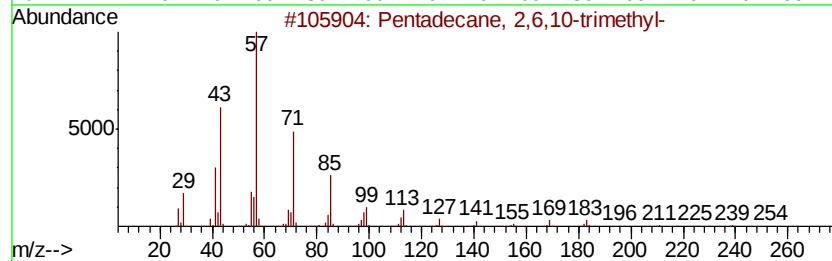
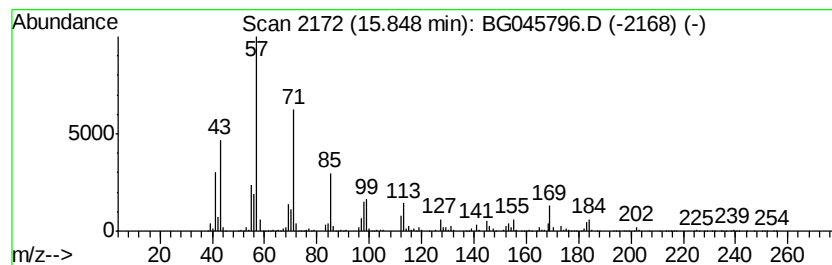
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane, 2,6,10-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.85	32.64 ng	6778770	Acenaphthene-d10	14.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	89
2		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	83
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	80
4		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	72
5		Tridecane	184	C13H28	000629-50-5	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
Sample : L3029-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WEST-17.5

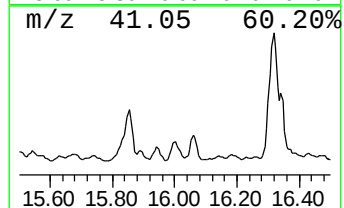
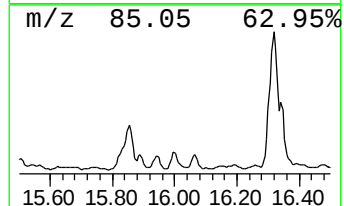
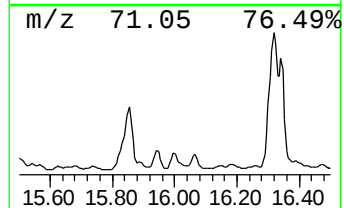
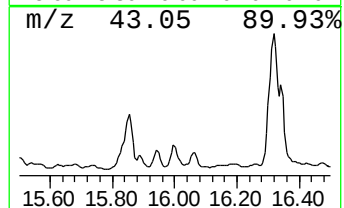
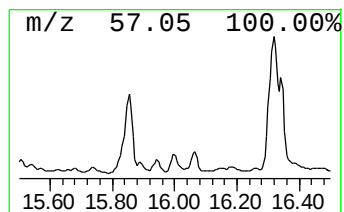
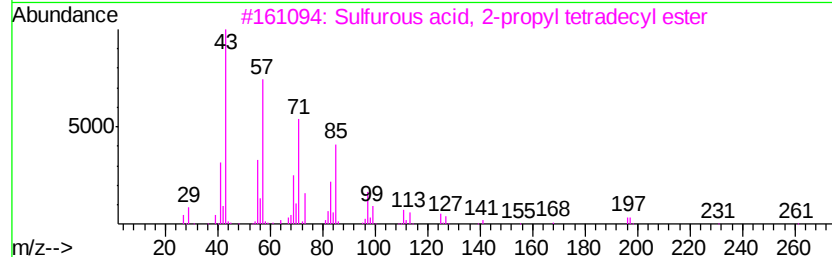
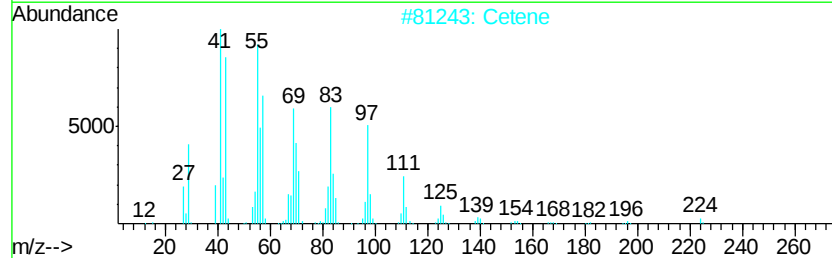
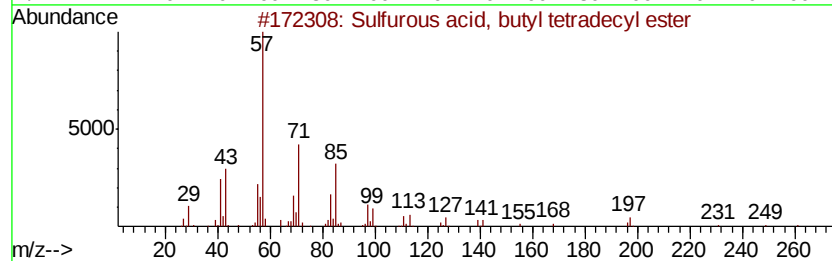
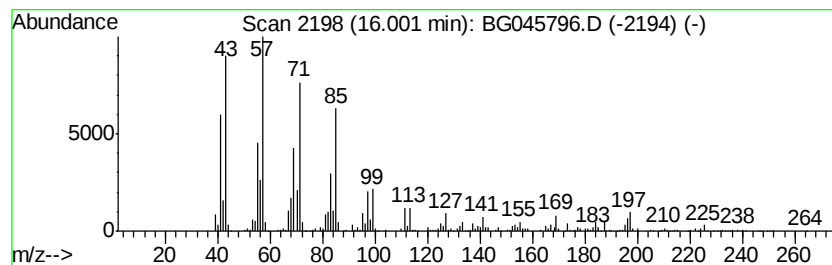
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Sulfurous acid, butyl tetra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.00	37.84 ng	2497500	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	91
2		Cetene	224	C16H32	000629-73-2	90
3		Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	87
4		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	87
5		Tetratetracontane	619	C44H90	007098-22-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
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ClientSampleId :
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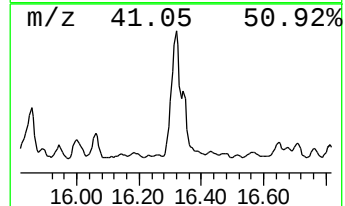
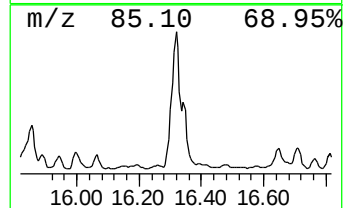
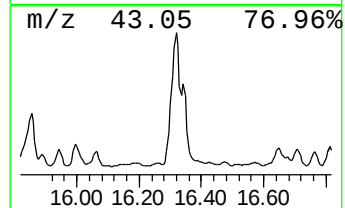
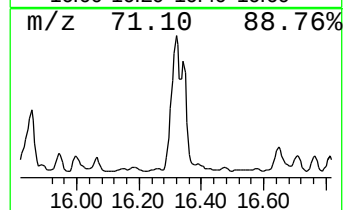
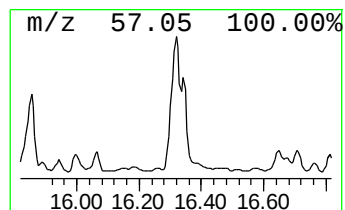
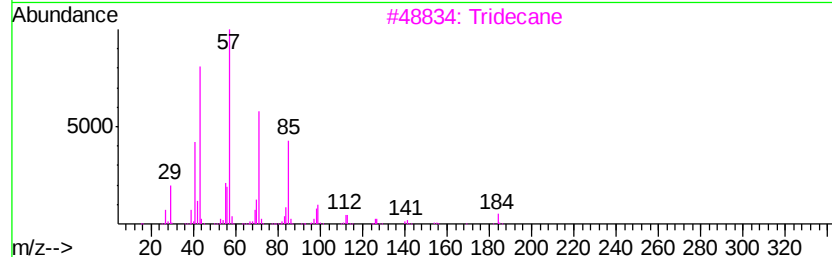
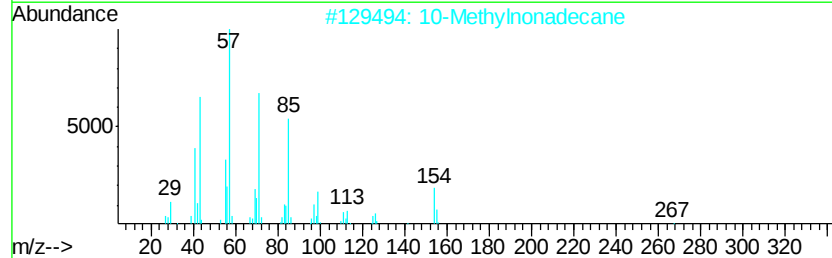
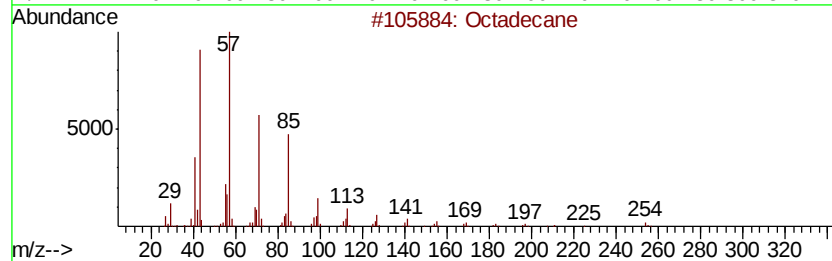
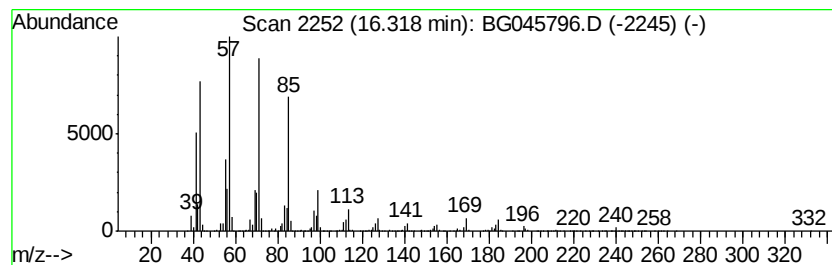
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.32	261.29 ng	17245200	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		10-Methylnonadecane	282	C20H42	056862-62-5	91
3		Tridecane	184	C13H28	000629-50-5	90
4		Nonadecane	268	C19H40	000629-92-5	90
5		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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Misc :
ALS Vial : 3 Sample Multiplier: 1

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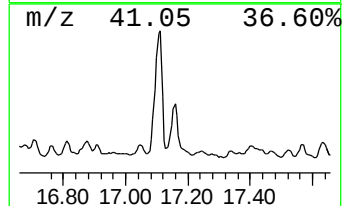
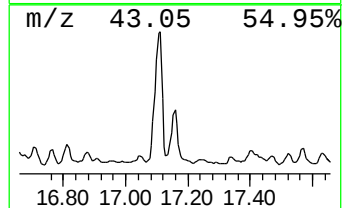
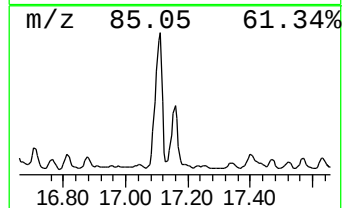
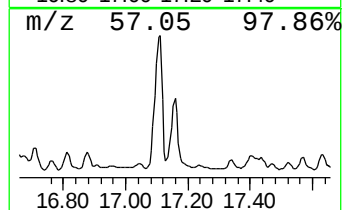
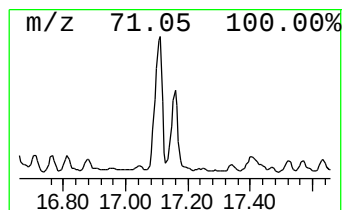
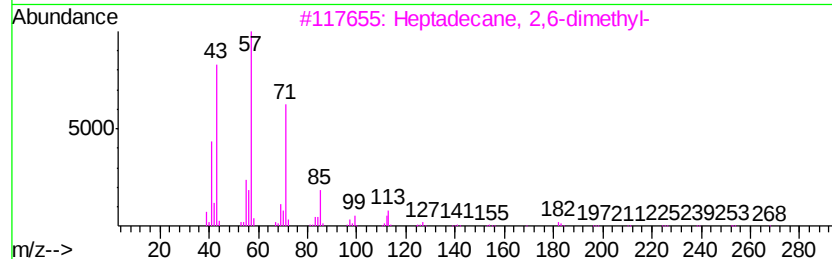
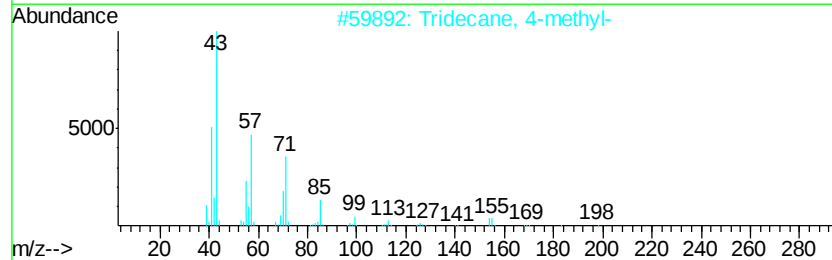
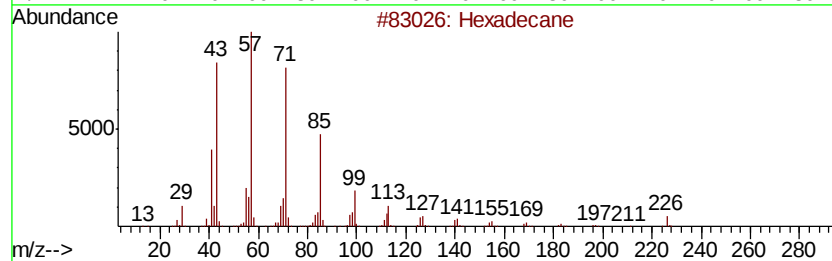
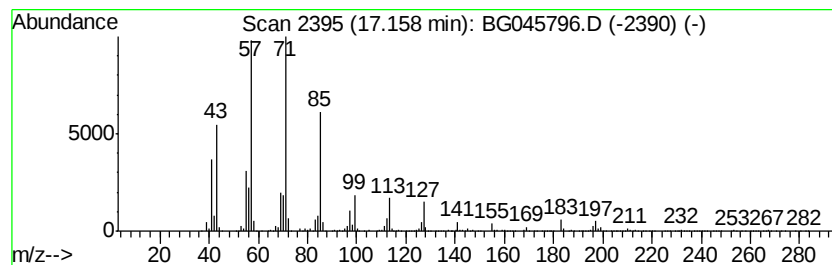
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.16	117.73 ng	7770210	Phenanthrene-d10	17.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Tridecane, 4-methyl-	198	C14H30	026730-12-1	86
3	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	83
4	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	83
5	2-Bromotetradecane	276	C14H29Br	074036-95-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
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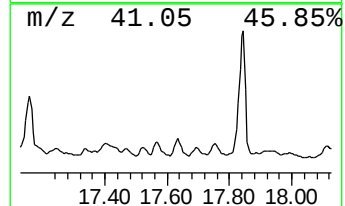
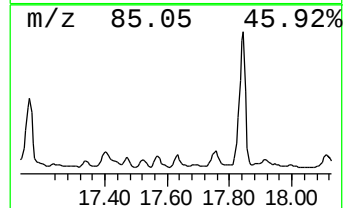
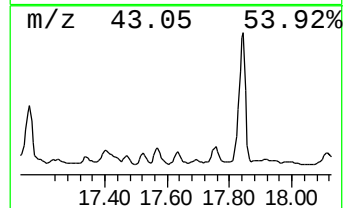
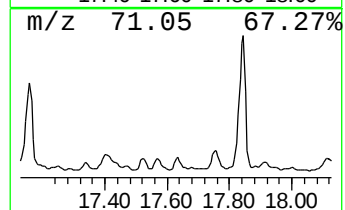
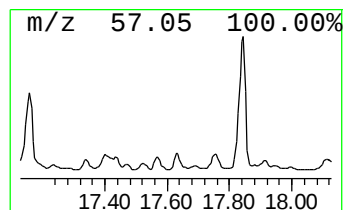
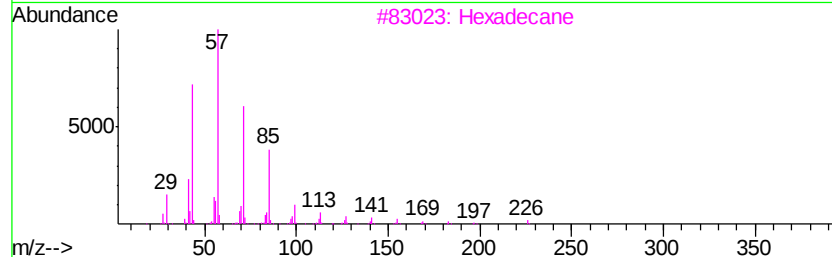
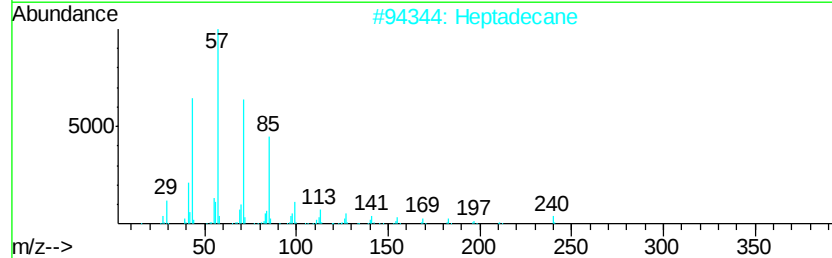
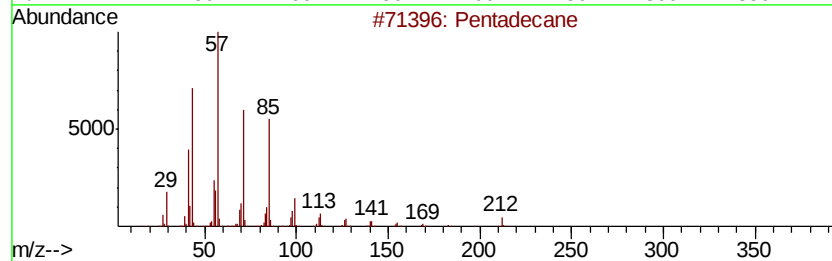
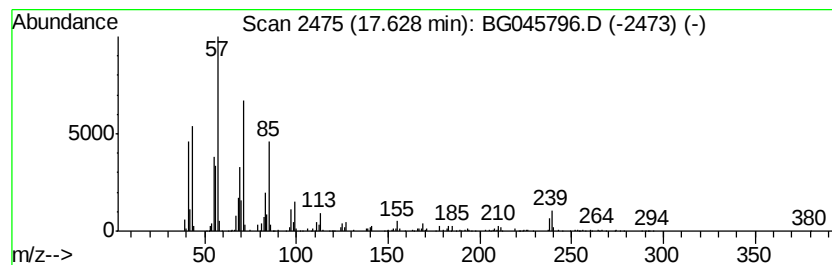
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Pentadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.63	23.02 ng	1519330	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	70
2		Heptadecane	240	C17H36	000629-78-7	70
3		Hexadecane	226	C16H34	000544-76-3	58
4		Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	49
5		Tetradecane, 3-methyl-	212	C15H32	018435-22-8	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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ClientSampleId :
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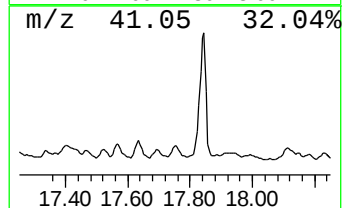
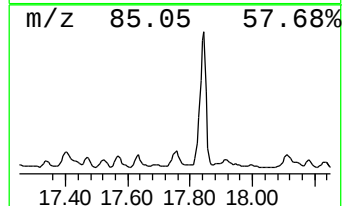
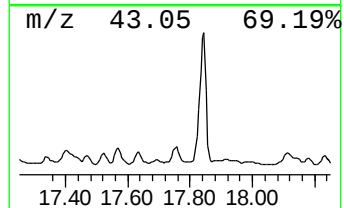
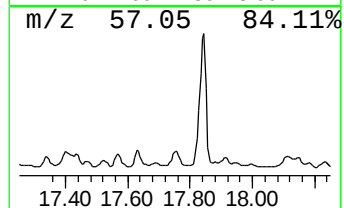
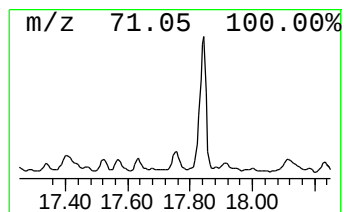
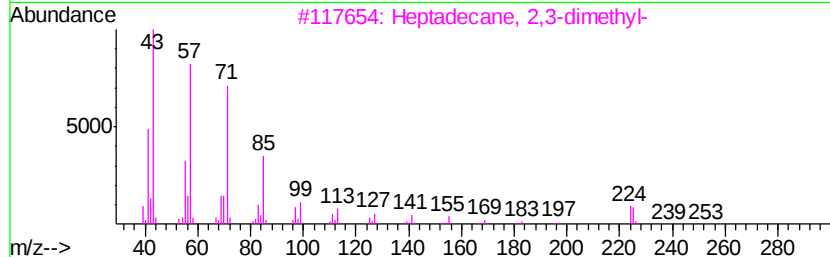
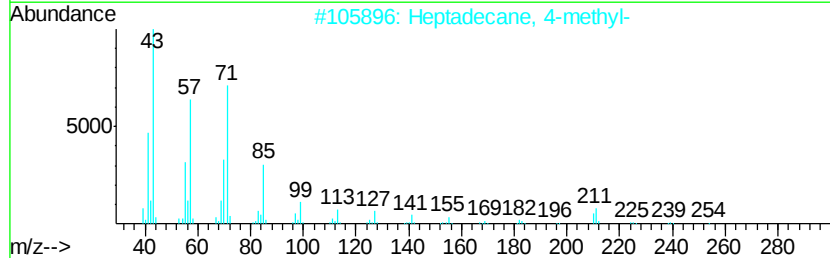
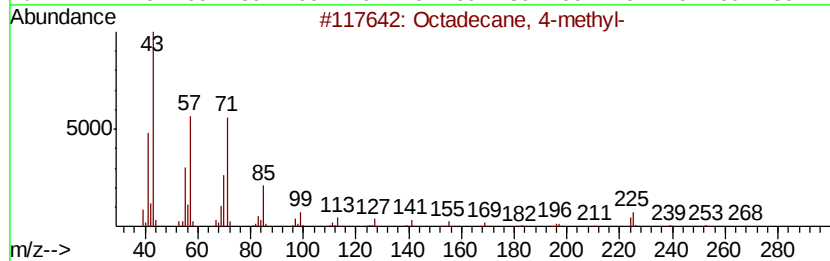
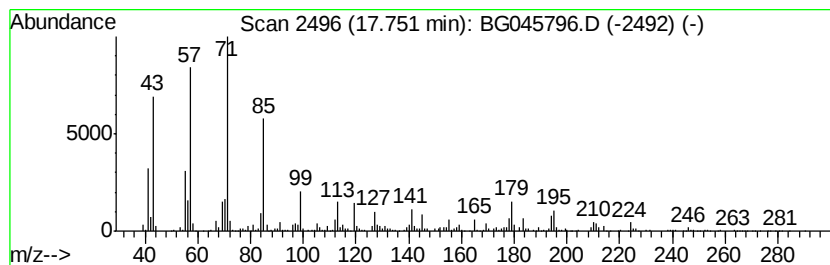
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecane, 4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.75	32.27 ng	2129900	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 4-methyl-	268	C19H40	010544-95-3	89
2		Heptadecane, 4-methyl-	254	C18H38	026429-11-8	81
3		Heptadecane, 2,3-dimethyl-	268	C19H40	061868-03-9	80
4		Octadecane, 2-methyl-	268	C19H40	001560-88-9	78
5		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	72



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Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
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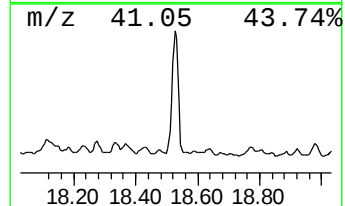
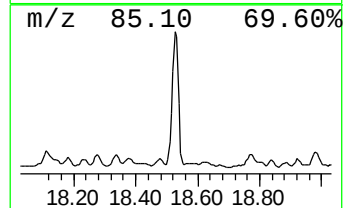
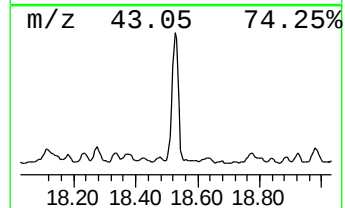
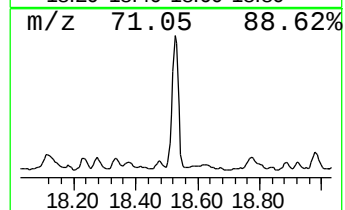
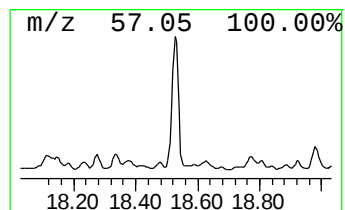
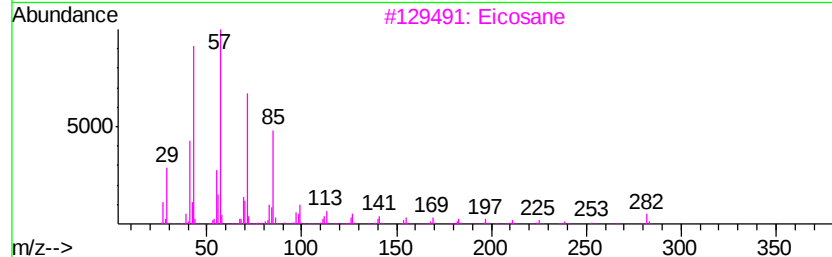
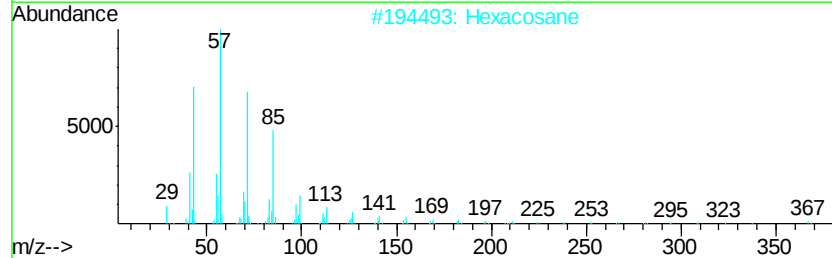
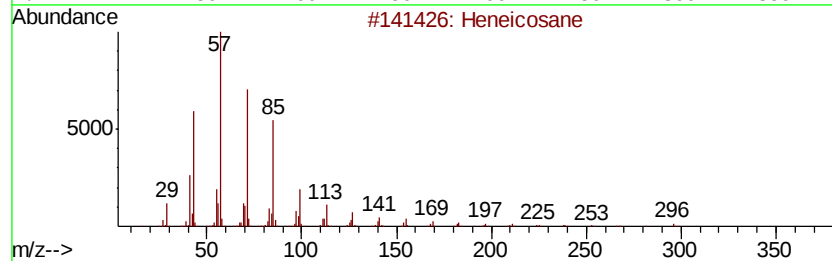
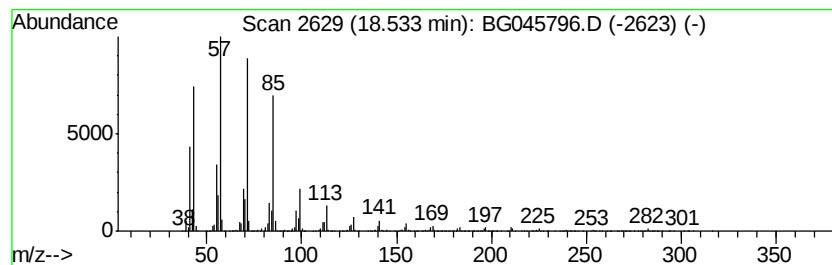
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.53	130.49 ng	8612460	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Eicosane	282	C20H42	000112-95-8	87
4		Nonadecane	268	C19H40	000629-92-5	83
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	83



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Misc :
ALS Vial : 3 Sample Multiplier: 1

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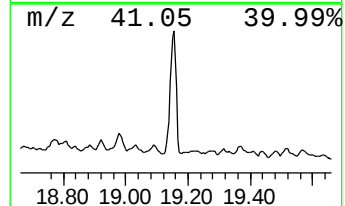
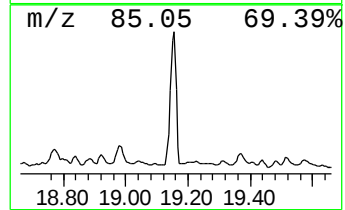
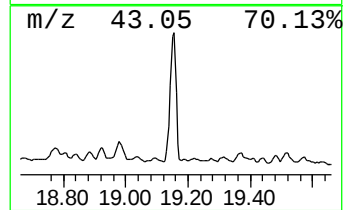
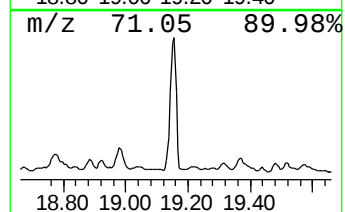
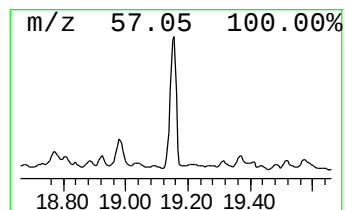
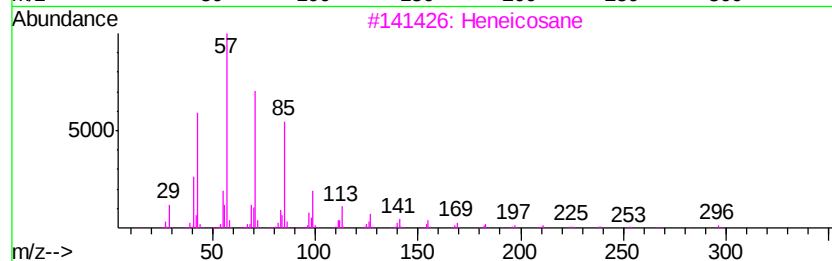
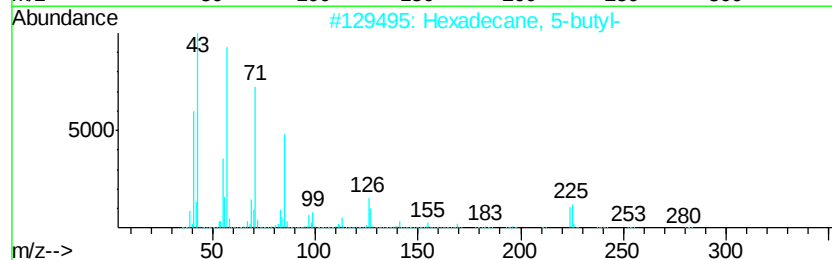
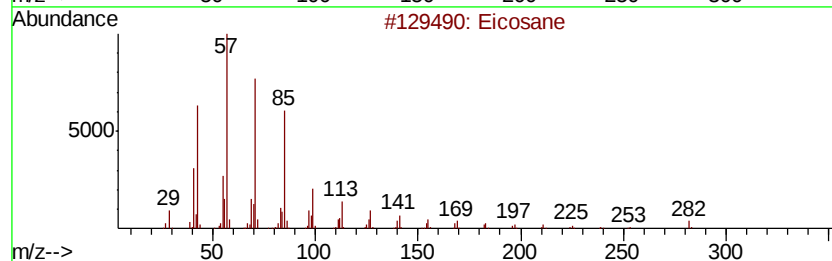
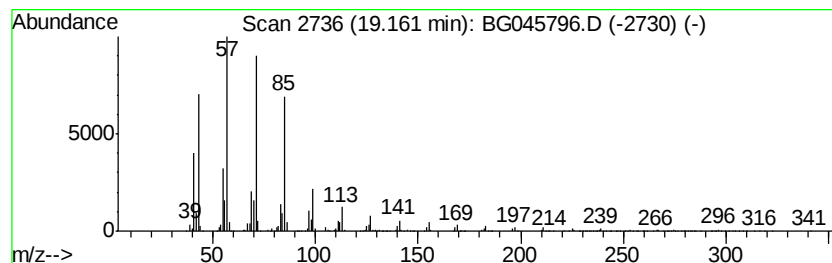
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Eicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	97.33 ng	6423590	Phenanthrene-d10	17.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Hexadecane, 5-butyl-	282	C20H42	006912-07-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
Data File : BG045796.D
Acq On : 18 Jun 2020 16:59
Operator : CG/JU
Sample : L3029-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WEST-17.5

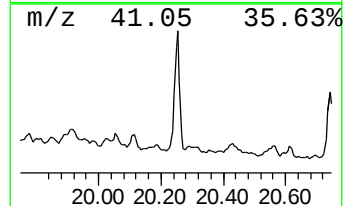
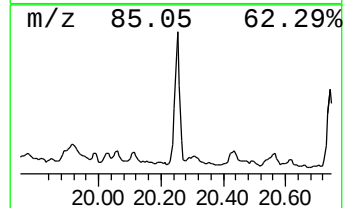
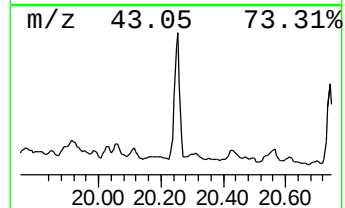
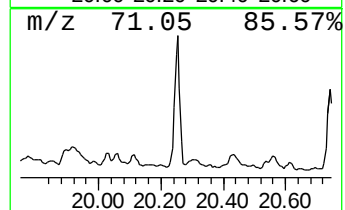
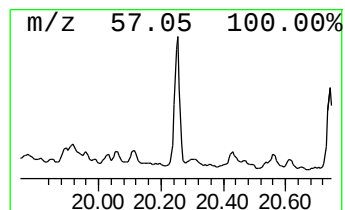
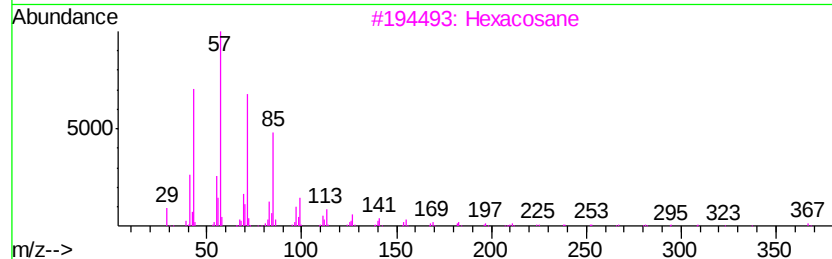
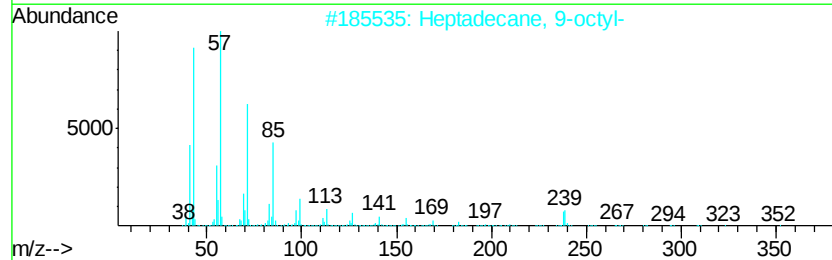
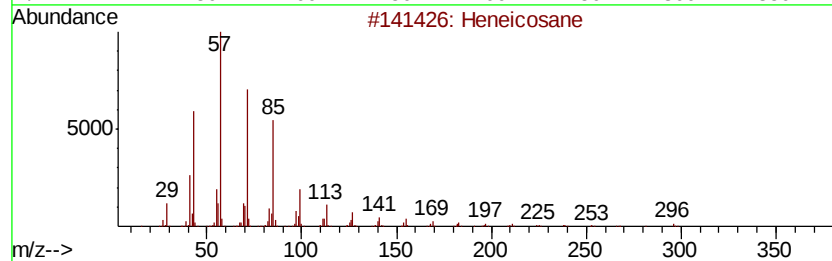
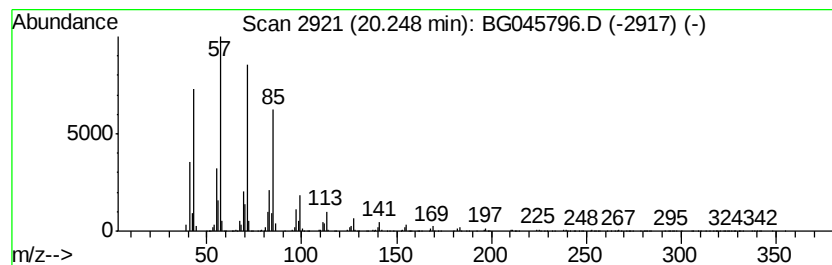
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Heptadecane, 9-octyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.25	30.06 ng	2937450	Chrysene-d12	21.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Eicosane	282	C20H42	000112-95-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG061820\
 Data File : BG045796.D
 Acq On : 18 Jun 2020 16:59
 Operator : CG/JU
 Sample : L3029-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WEST-17.5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG061520.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
Nonane	5.94	22.4	ng	2092360	1	7.92	1867800	20.0
Decane	7.56	64.0	ng	5978950	1	7.92	1867800	20.0
Benzene, 1-methyl...	8.47	29.1	ng	2715750	1	7.92	1867800	20.0
Benzene, 1,2-diet...	8.57	44.2	ng	4125910	1	7.92	1867800	20.0
Octadecane, 1-chl...	9.17	73.7	ng	6880300	1	7.92	1867800	20.0
Benzene, 4-(chlor...	9.29	37.0	ng	3452040	1	7.92	1867800	20.0
Naphthalene, 1,2,...	13.53	24.4	ng	5058730	3	14.59	4154190	20.0
Nonadecane	14.51	73.3	ng	15216400	3	14.59	4154190	20.0
10-Methylnonadecane	15.46	67.6	ng	14047600	3	14.59	4154190	20.0
Pentadecane, 2,6,...	15.85	32.6	ng	6778770	3	14.59	4154190	20.0
Sulfurous acid, b...	16.00	37.8	ng	2497500	4	17.33	1320010	20.0
Octadecane	16.32	261.3	ng	17245200	4	17.33	1320010	20.0
Hexadecane	17.16	117.7	ng	7770210	4	17.33	1320010	20.0
Pentadecane	17.63	23.0	ng	1519330	4	17.33	1320010	20.0
Octadecane, 4-met...	17.75	32.3	ng	2129900	4	17.33	1320010	20.0
Heneicosane	18.53	130.5	ng	8612460	4	17.33	1320010	20.0
Eicosane	19.16	97.3	ng	6423590	4	17.33	1320010	20.0
Heptadecane, 9-oc...	20.25	30.1	ng	2937450	5	21.58	1954690	20.0