

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
 Data File : BG064162.D
 Acq On : 2 Apr 2025 21:04
 Operator : RC/JU
 Sample : Q1692-01 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-08-04012025

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_G\Methods\8270-BG030525.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064162.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.992	318	324	331	rBV2	38169	61433	9.73%	0.857%
2	5.450	395	402	414	rVB	138091	241382	38.25%	3.368%
3	6.067	502	507	514	rVB4	5247	10542	1.67%	0.147%
4	7.025	665	670	679	rVB	146380	256191	40.59%	3.574%
5	7.865	804	813	820	rBV	161907	275461	43.65%	3.843%
6	9.011	1002	1008	1016	rVB2	92897	159645	25.30%	2.227%
7	9.105	1018	1024	1030	rVB2	8196	14623	2.32%	0.204%
8	9.587	1100	1106	1110	rVB6	3861	7030	1.11%	0.098%
9	10.650	1278	1287	1293	rBV	218914	419503	66.47%	5.853%
10	10.821	1314	1316	1321	rVB4	7311	11616	1.84%	0.162%
11	11.326	1397	1402	1403	rBV4	4009	6866	1.09%	0.096%
12	11.502	1428	1432	1436	rBV4	6413	9916	1.57%	0.138%
13	11.573	1438	1444	1451	rBV5	9016	20138	3.19%	0.281%
14	11.684	1458	1463	1468	rBV3	11592	22315	3.54%	0.311%
15	11.843	1485	1490	1498	rVB4	11851	21275	3.37%	0.297%
16	12.084	1524	1531	1537	rBV3	35436	61110	9.68%	0.853%
17	12.195	1546	1550	1553	rBV5	6102	8039	1.27%	0.112%
18	12.307	1565	1569	1577	rVB7	10512	18039	2.86%	0.252%
19	12.530	1602	1607	1609	rBV5	7339	11875	1.88%	0.166%
20	12.560	1609	1612	1618	rVB4	14474	24306	3.85%	0.339%
21	12.613	1618	1621	1624	rBV4	4590	6732	1.07%	0.094%
22	12.759	1643	1646	1651	rVB6	6613	10976	1.74%	0.153%
23	12.824	1651	1657	1660	rBV5	7845	12509	1.98%	0.175%
24	12.900	1666	1670	1674	rVB5	7209	11206	1.78%	0.156%
25	12.983	1679	1684	1689	rBV6	12165	20815	3.30%	0.290%
26	13.036	1689	1693	1699	rVV4	19018	27158	4.30%	0.379%
27	13.106	1699	1705	1713	rVV	227012	365205	57.87%	5.095%
28	13.194	1717	1720	1723	rVB5	5602	7621	1.21%	0.106%
29	13.324	1737	1742	1749	rBV	56748	84561	13.40%	1.180%
30	13.435	1756	1761	1766	rBV7	10198	20132	3.19%	0.281%
31	13.494	1766	1771	1773	rVV5	5875	7416	1.18%	0.103%
32	13.570	1782	1784	1789	rVB5	7370	10513	1.67%	0.147%
33	13.653	1793	1798	1799	rBV4	7153	8994	1.43%	0.125%
34	13.764	1814	1817	1820	rBV5	7744	12568	1.99%	0.175%
35	13.852	1829	1832	1837	rVB7	11201	14876	2.36%	0.208%
36	13.958	1846	1850	1852	rBV4	9418	13333	2.11%	0.186%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.982	1852	1854	1858	rVV5	18750	26654	4.22%	0.372%
38	14.017	1858	1860	1867	rVB6	10333	16316	2.59%	0.228%
39	14.099	1873	1874	1878	rVB4	11313	9566	1.52%	0.133%
40	14.146	1878	1882	1883	rBV4	6490	6314	1.00%	0.088%
41	14.199	1887	1891	1895	rVB5	5709	9175	1.45%	0.128%
42	14.258	1897	1901	1904	rVB6	5707	7044	1.12%	0.098%
43	14.328	1912	1913	1919	rVB5	7101	9044	1.43%	0.126%
44	14.399	1919	1925	1929	rBV	85792	109625	17.37%	1.529%
45	14.487	1929	1940	1946	rBV	340174	548995	86.99%	7.659%
46	14.551	1948	1951	1956	rVB5	13277	19774	3.13%	0.276%
47	14.616	1961	1962	1968	rVB6	6794	8751	1.39%	0.122%
48	14.845	1998	2001	2002	rBV3	8032	8771	1.39%	0.122%
49	14.898	2008	2010	2015	rVB5	10776	13341	2.11%	0.186%
50	14.957	2018	2020	2023	rVB3	7172	8047	1.28%	0.112%
51	15.022	2025	2031	2039	rVB7	19828	38234	6.06%	0.533%
52	15.086	2039	2042	2045	rBV4	11115	15562	2.47%	0.217%
53	15.274	2072	2074	2079	rBV6	4911	6707	1.06%	0.094%
54	15.351	2083	2087	2092	rVB	87989	110327	17.48%	1.539%
55	15.450	2101	2104	2110	rVB7	7912	12668	2.01%	0.177%
56	15.533	2113	2118	2121	rBV7	14409	25301	4.01%	0.353%
57	15.574	2122	2125	2132	rVV9	4464	9372	1.48%	0.131%
58	15.697	2144	2146	2149	rBV3	8398	8952	1.42%	0.125%
59	15.756	2150	2156	2160	rVV2	29626	53137	8.42%	0.741%
60	15.838	2166	2170	2177	rVB5	33520	54178	8.58%	0.756%
61	15.903	2177	2181	2187	rBV6	13321	28466	4.51%	0.397%
62	15.967	2187	2192	2201	rVB2	200956	305110	48.34%	4.257%
63	16.044	2201	2205	2208	rBV6	7940	11381	1.80%	0.159%
64	16.167	2224	2226	2229	rVB4	7252	6415	1.02%	0.089%
65	16.208	2229	2233	2236	rBV	96341	129613	20.54%	1.808%
66	16.320	2250	2252	2260	rBV9	6141	11873	1.88%	0.166%
67	16.608	2300	2301	2306	rVB4	11471	10778	1.71%	0.150%
68	16.667	2309	2311	2316	rVB5	6692	9034	1.43%	0.126%
69	16.778	2328	2330	2334	rBV3	7005	9215	1.46%	0.129%
70	17.002	2364	2368	2372	rBV	85766	107668	17.06%	1.502%
71	17.049	2373	2376	2381	rVB3	32586	51324	8.13%	0.716%
72	17.225	2399	2406	2411	rBV2	409443	631117	100.00%	8.805%
73	17.266	2411	2413	2417	rVB	68104	83303	13.20%	1.162%
74	17.354	2425	2428	2434	rVB3	19505	31338	4.97%	0.437%
75	17.430	2438	2441	2444	rVV4	9841	12595	2.00%	0.176%
76	17.466	2445	2447	2452	rVB5	12244	13610	2.16%	0.190%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	17.542	2457	2460	2463	rVB3	11423	13537	2.14%	0.189%
78	17.660	2477	2480	2486	rVB7	11904	16897	2.68%	0.236%
79	17.736	2489	2493	2497	rBV2	82834	99948	15.84%	1.394%
80	17.801	2502	2504	2509	rVV6	6372	7628	1.21%	0.106%
81	17.953	2525	2530	2531	rBV5	10342	15446	2.45%	0.215%
82	18.024	2537	2542	2546	rBV8	9869	21532	3.41%	0.300%
83	18.094	2552	2554	2556	rBV2	14399	15227	2.41%	0.212%
84	18.294	2583	2588	2591	rBV5	21311	40518	6.42%	0.565%
85	18.429	2607	2611	2615	rVB	72574	88215	13.98%	1.231%
86	18.717	2658	2660	2661	rBV2	9141	6782	1.07%	0.095%
87	19.058	2715	2718	2722	rBV	66263	80034	12.68%	1.117%
88	19.281	2750	2756	2761	rBV	83721	131977	20.91%	1.841%
89	19.640	2813	2817	2821	rVB2	128790	161920	25.66%	2.259%
90	19.845	2847	2852	2858	rBV	375934	490150	77.66%	6.838%
91	21.455	3121	3126	3132	rBV	392335	607868	96.32%	8.481%
92	24.475	3634	3640	3649	rVB2	200597	485495	76.93%	6.773%

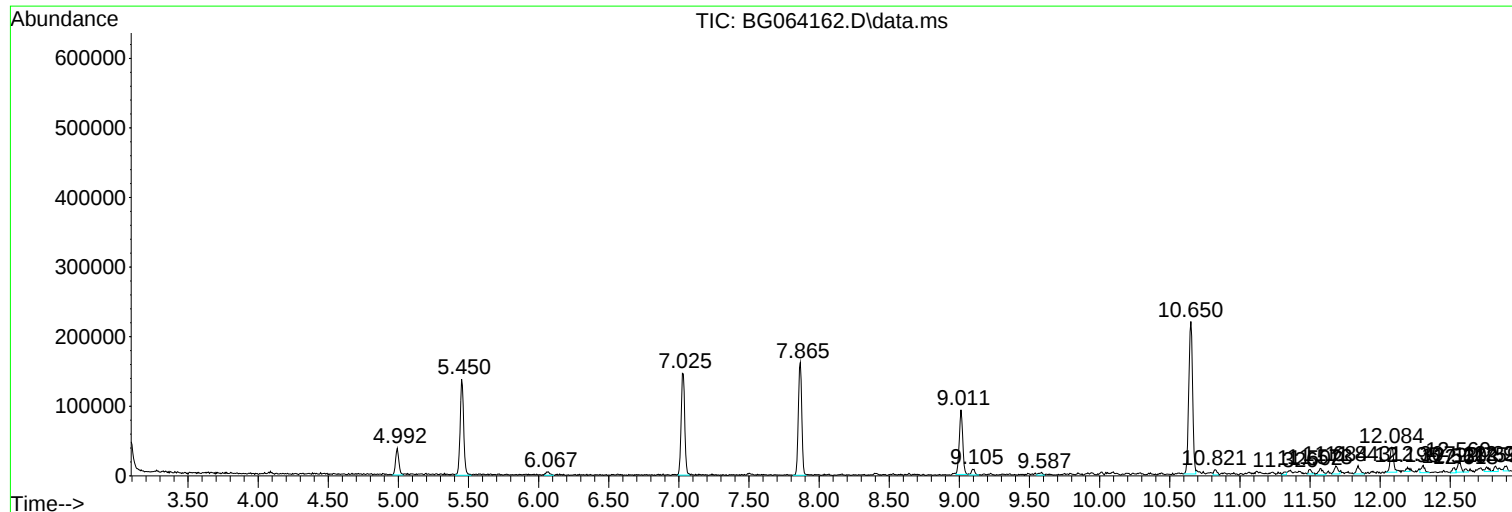
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Instrument :
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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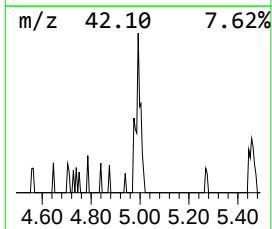
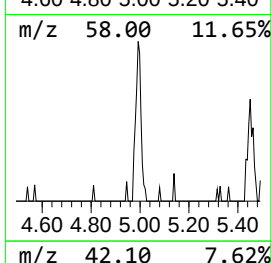
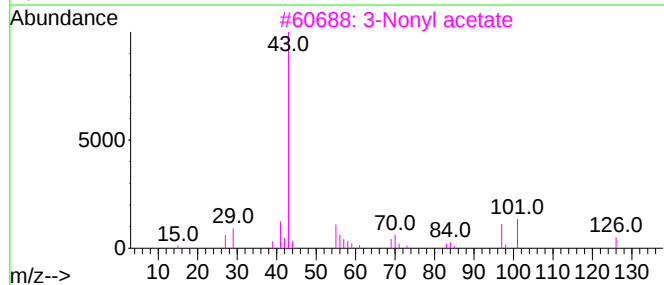
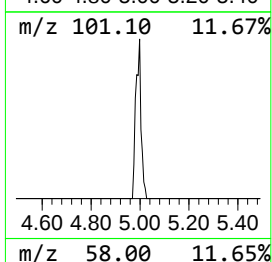
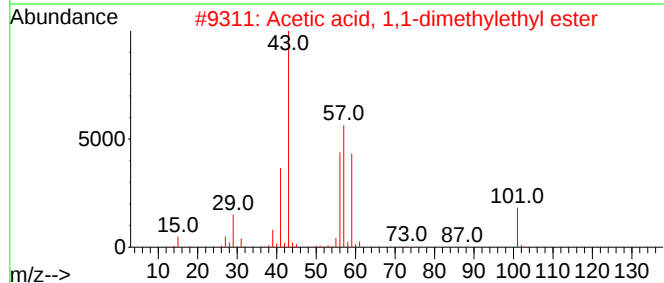
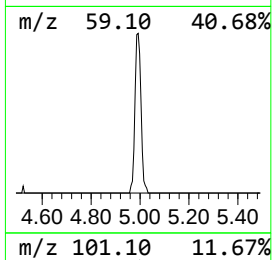
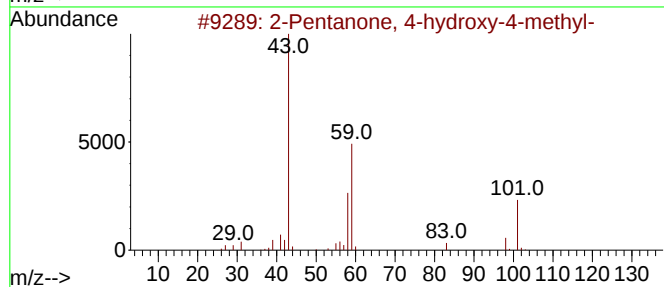
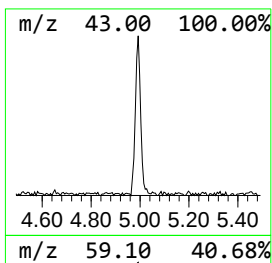
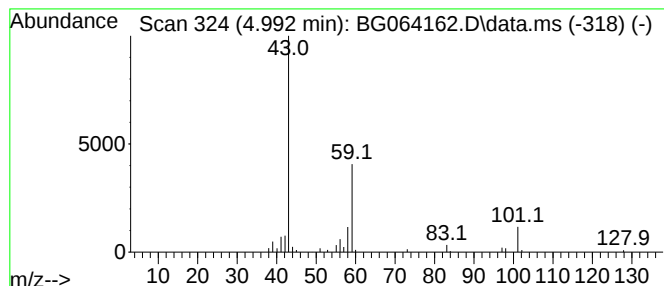
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.992	4.46 ng	61433	1,4-Dichlorobenzene-d4	7.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	33		
3	3-Nonyl acetate	186	C11H22O2	1010429-16-2	17		
4	2-Pentanone, 5-(acetyloxy)-	144	C7H12O3	005185-97-7	17		
5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9		



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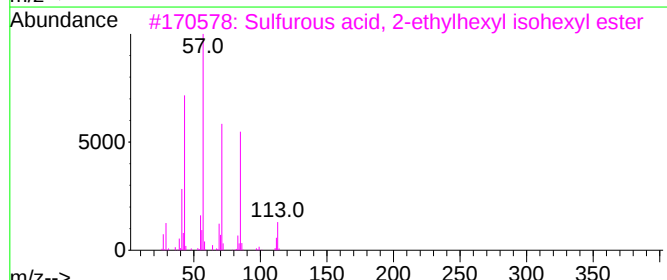
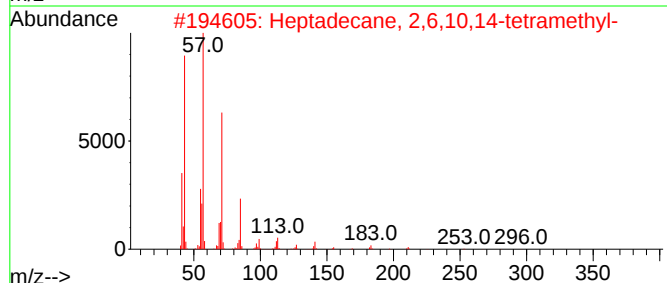
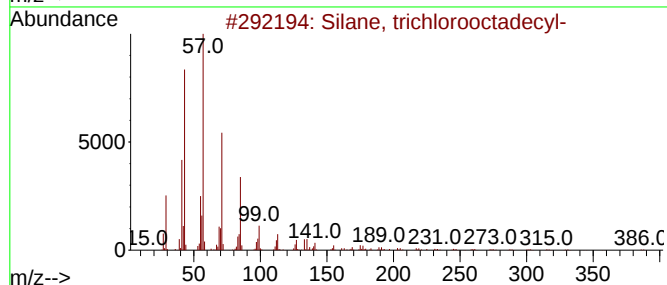
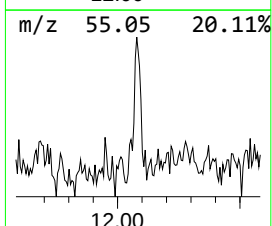
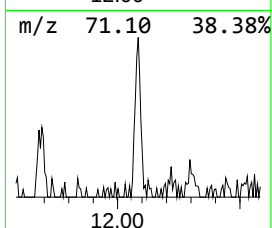
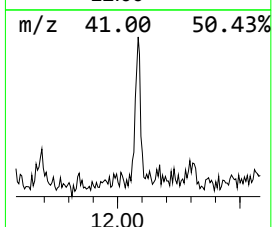
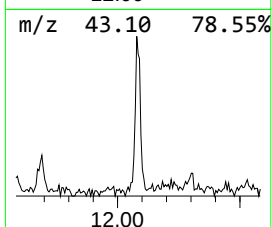
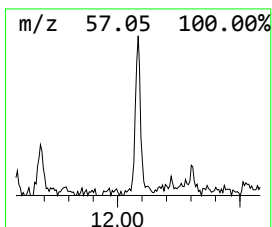
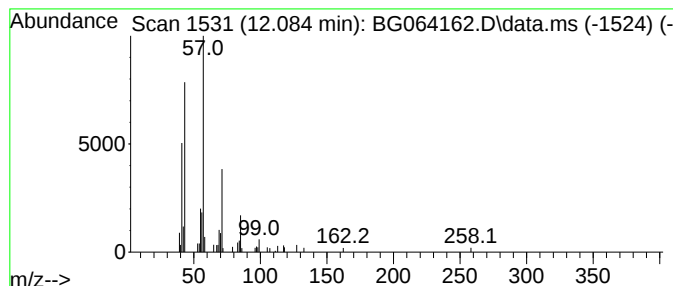
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Silane, trichlorooctadecyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.084	2.91 ng	61110	Naphthalene-d8	10.650

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	83
2		Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	83
3		Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	72
4		Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	72
5		Tridecane	184	C13H28	000629-50-5	72



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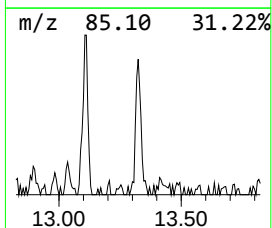
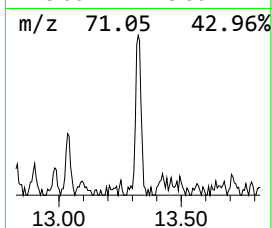
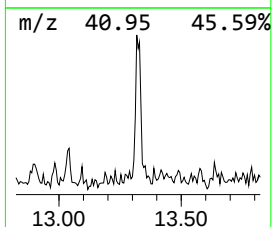
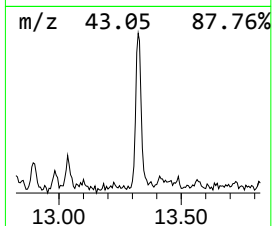
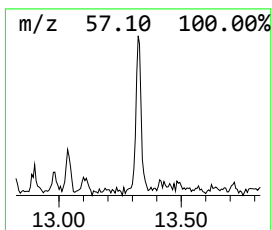
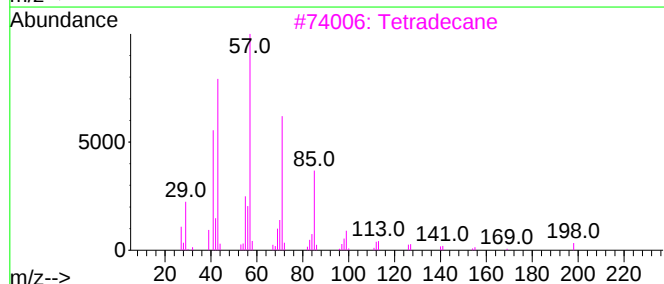
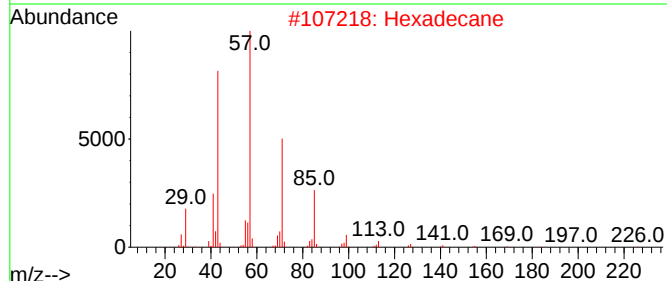
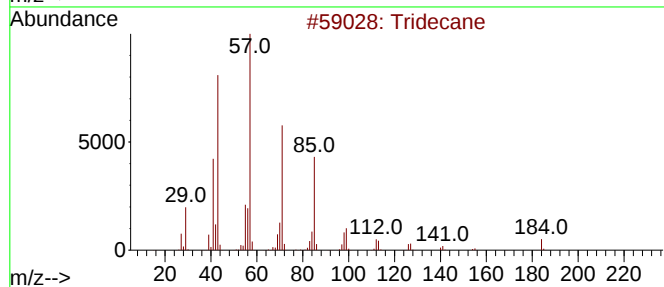
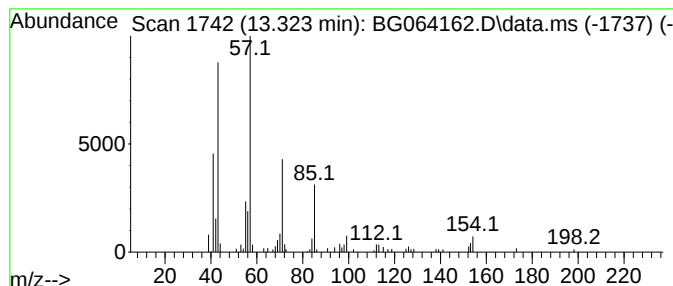
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Tridecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.323	3.08 ng	84561	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	80
2			Hexadecane	226	C16H34	000544-76-3	80
3			Tetradecane	198	C14H30	000629-59-4	74
4			Pentadecane	212	C15H32	000629-62-9	72
5			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	64



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ClientSampleId :
NB-08-04012025

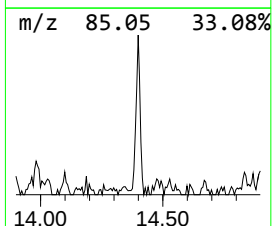
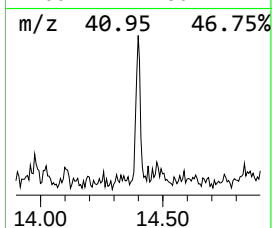
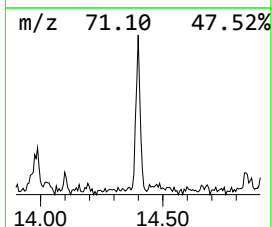
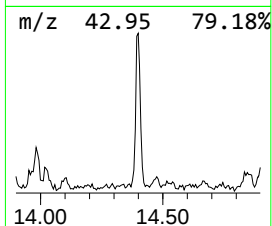
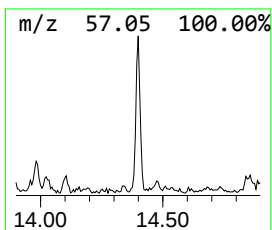
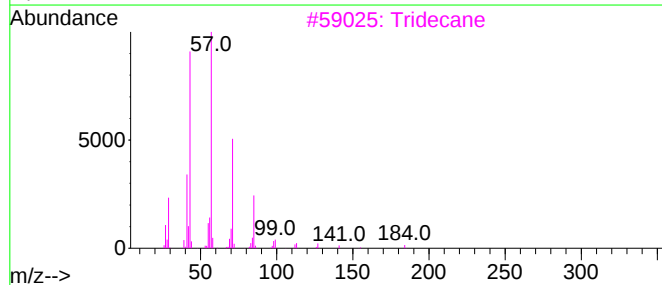
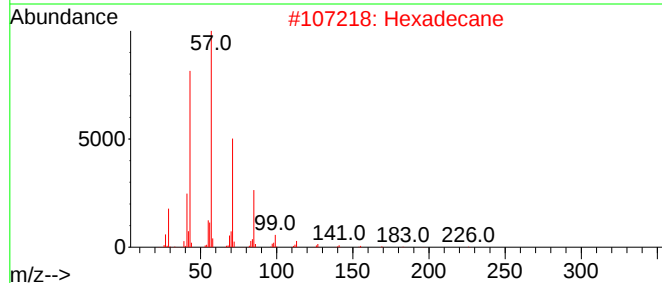
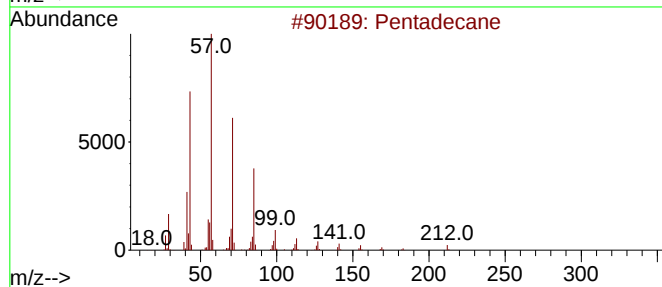
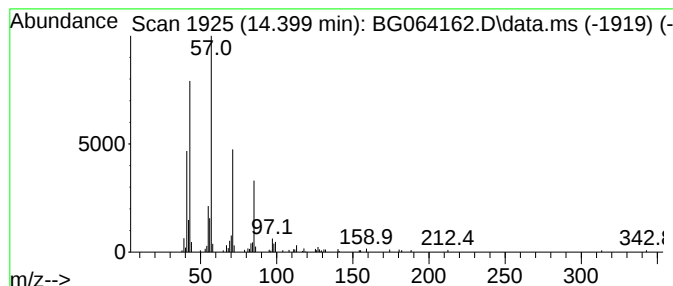
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.399	3.99 ng	109625	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	91
2			Hexadecane	226	C16H34	000544-76-3	90
3			Tridecane	184	C13H28	000629-50-5	83
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
5			1-Iodoundecane	282	C11H23I	004282-44-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-04012025

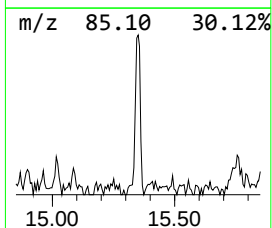
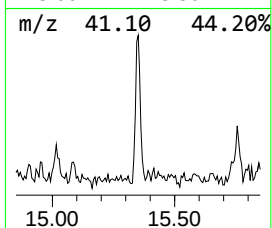
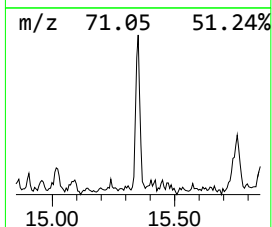
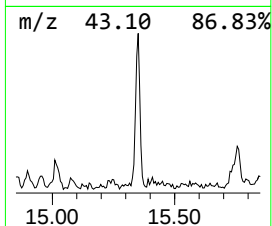
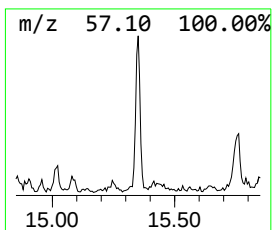
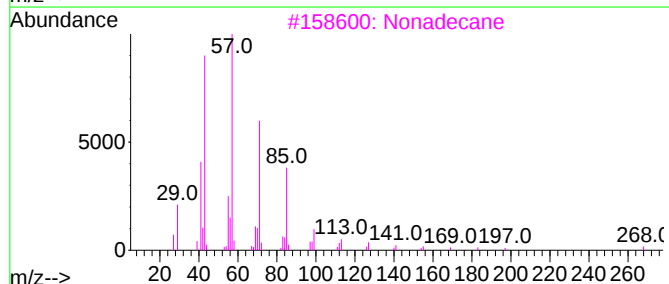
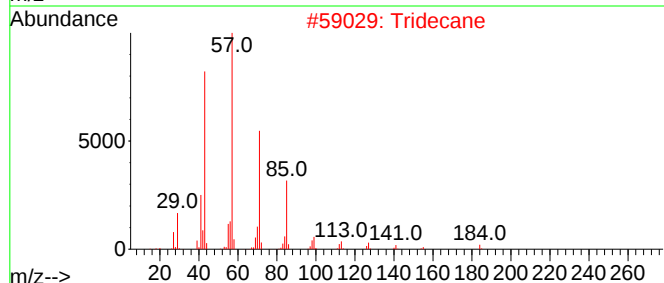
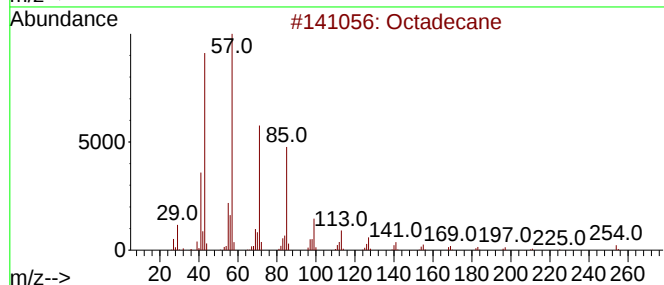
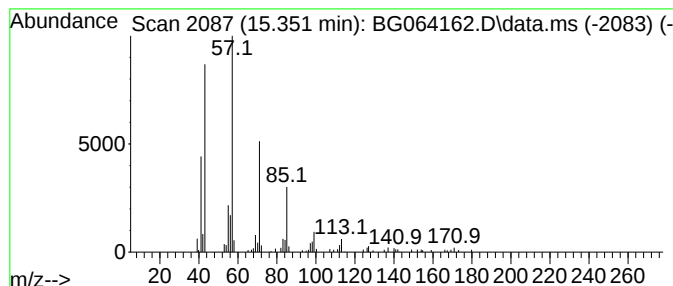
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.351	4.02 ng	110327	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	90
2			Tridecane	184	C13H28	000629-50-5	90
3			Nonadecane	268	C19H40	000629-92-5	90
4			Hexadecane	226	C16H34	000544-76-3	90
5			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-04012025

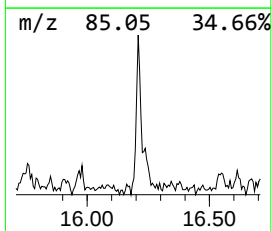
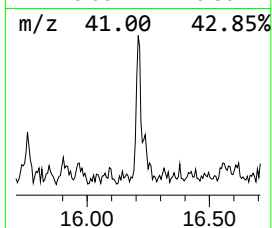
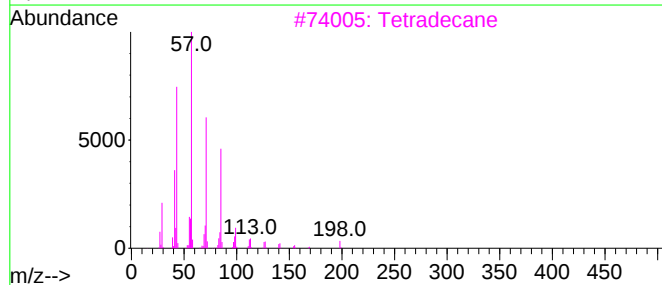
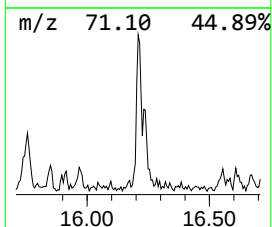
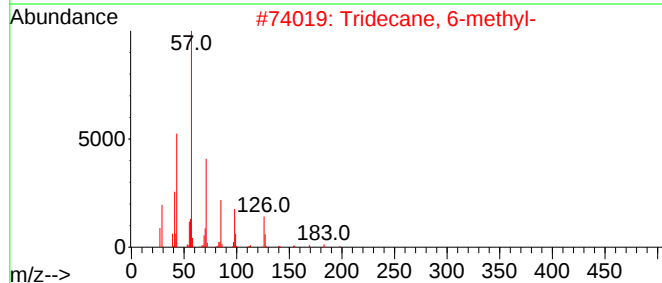
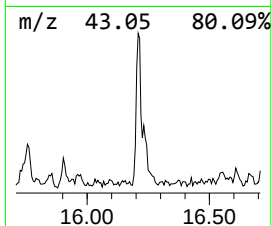
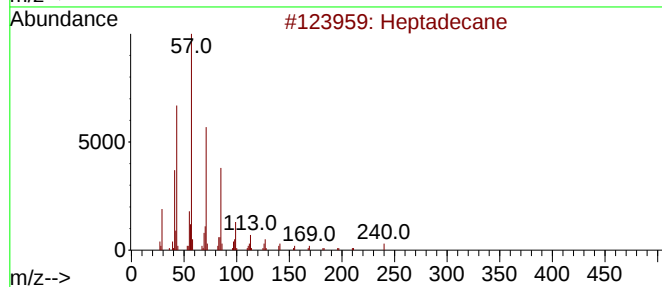
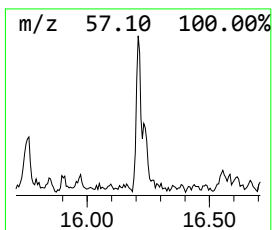
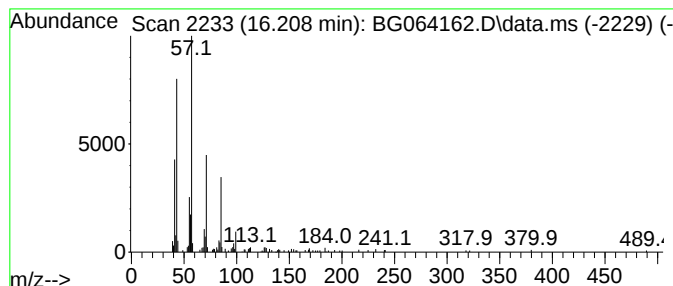
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptadecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.208	4.11 ng	129613	Phenanthrene-d10	17.225

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	83
2			Tridecane, 6-methyl-	198	C14H30	013287-21-3	83
3			Tetradecane	198	C14H30	000629-59-4	81
4			Tridecane, 7-propyl-	226	C16H34	055045-09-5	72
5			Dodecane, 2-methyl-	184	C13H28	001560-97-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-04012025

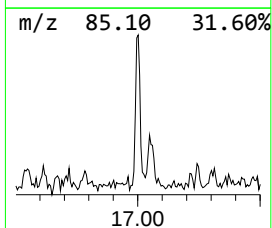
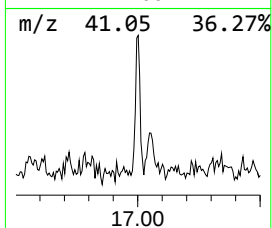
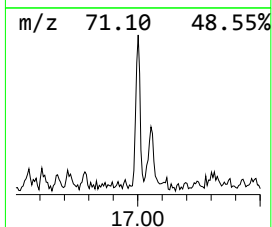
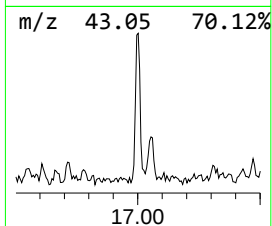
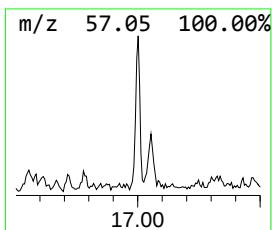
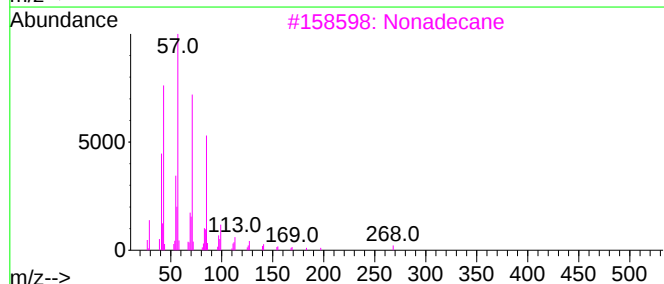
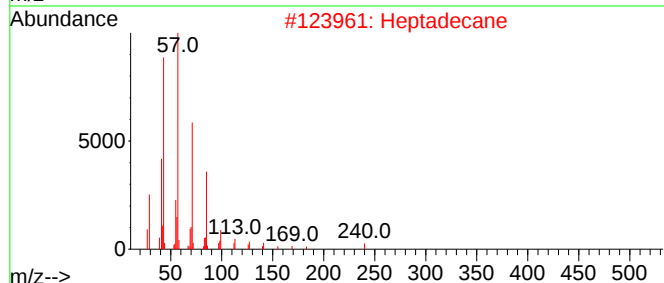
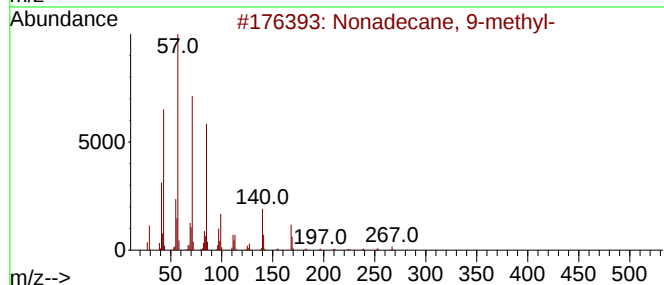
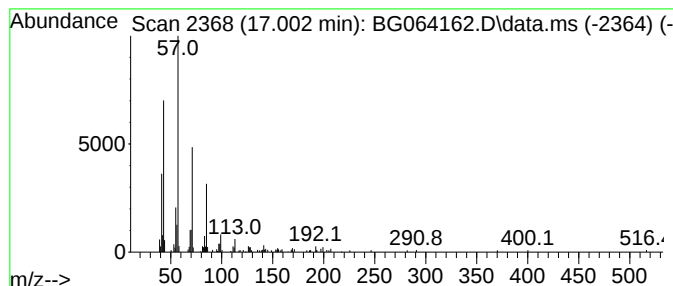
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Nonadecane, 9-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.002	3.41 ng	107668	Phenanthrene-d10	17.225

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	87
2			Heptadecane	240	C17H36	000629-78-7	80
3			Nonadecane	268	C19H40	000629-92-5	74
4			Eicosane	282	C20H42	000112-95-8	74
5			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NB-08-04012025

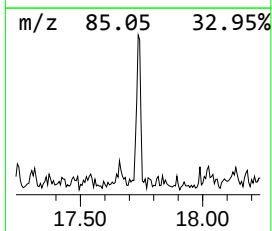
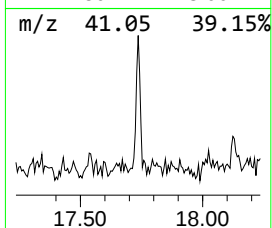
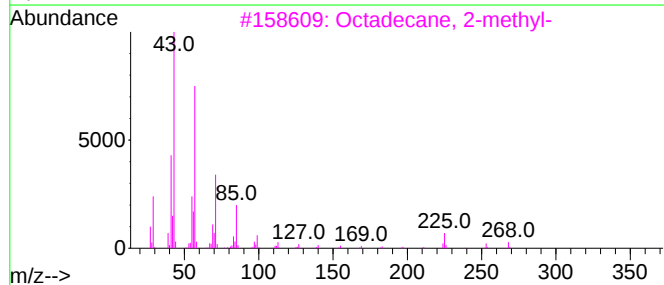
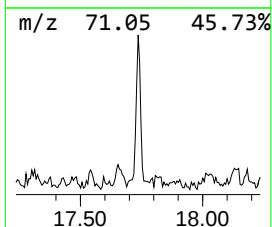
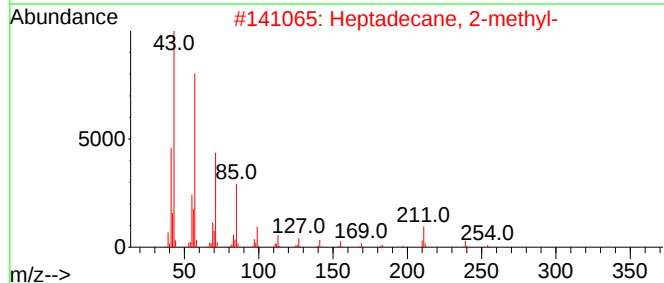
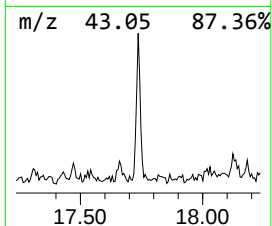
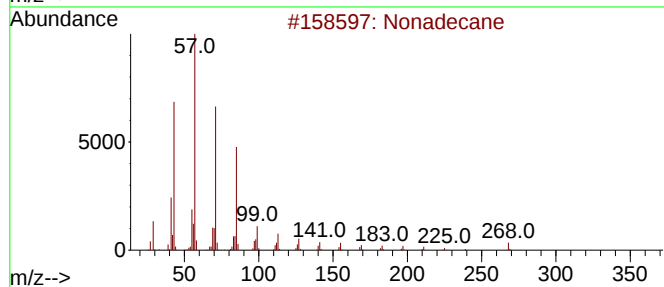
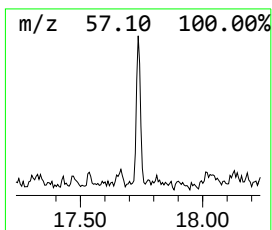
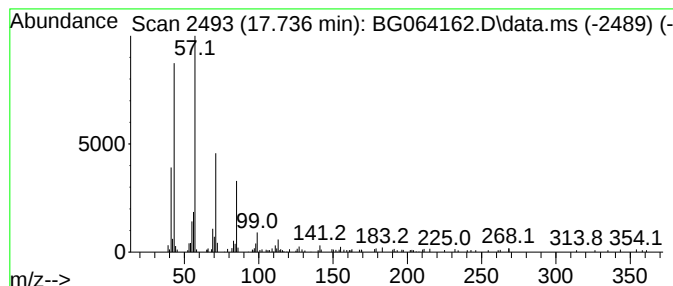
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.736	3.17 ng	99948	Phenanthrene-d10	17.225

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	87
2			Heptadecane, 2-methyl-	254	C18H38	001560-89-0	81
3			Octadecane, 2-methyl-	268	C19H40	001560-88-9	80
4			Eicosane	282	C20H42	000112-95-8	72
5			Octadecane	254	C18H38	000593-45-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NB-08-04012025

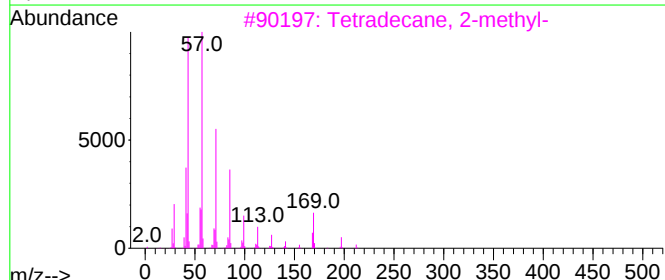
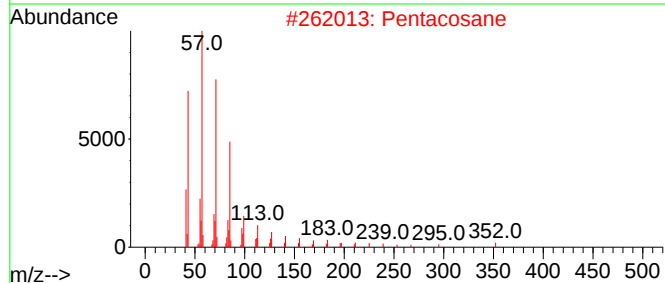
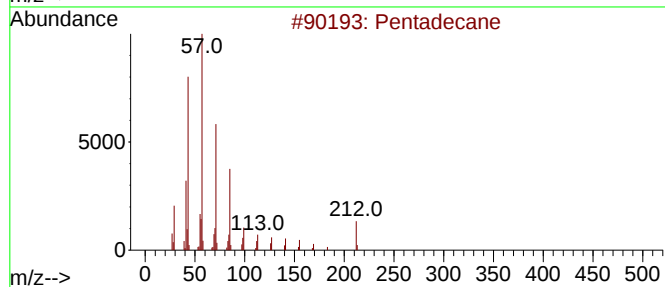
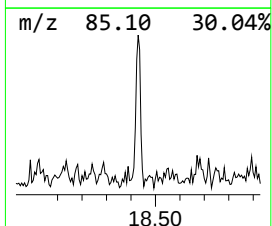
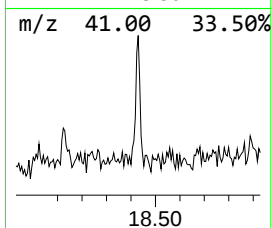
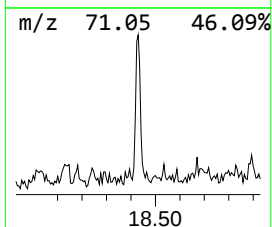
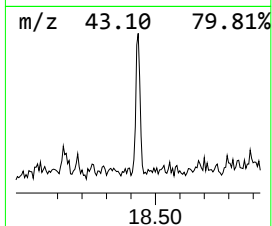
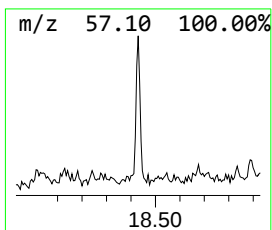
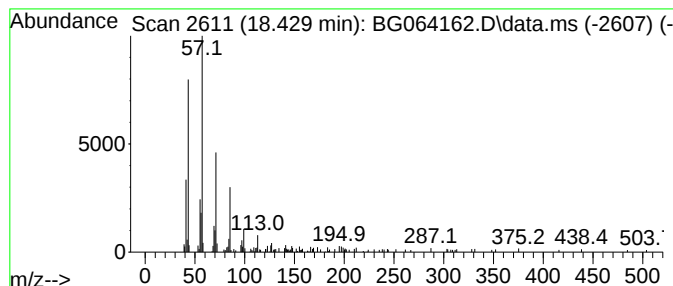
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Pentacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.429	2.80 ng	88215	Phenanthrene-d10	17.225

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	91
2			Pentacosane	352	C25H52	000629-99-2	89
3			Tetradecane, 2-methyl-	212	C15H32	001560-95-8	87
4			Nonadecane, 2-methyl-	282	C20H42	001560-86-7	87
5			Octadecane, 2-methyl-	268	C19H40	001560-88-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
NB-08-04012025

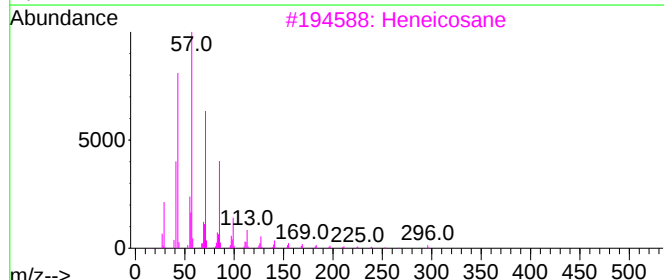
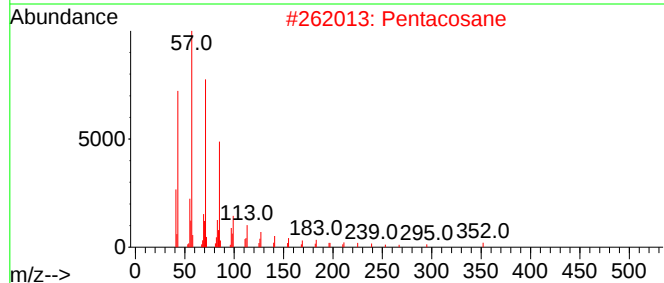
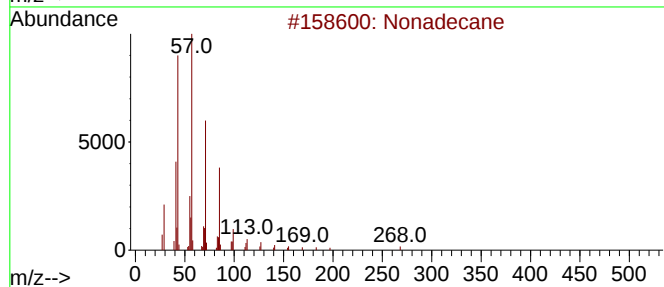
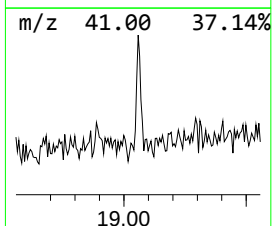
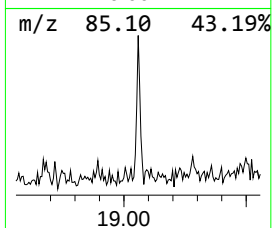
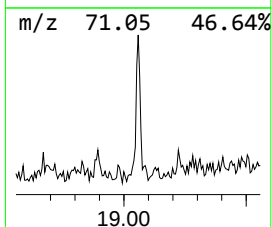
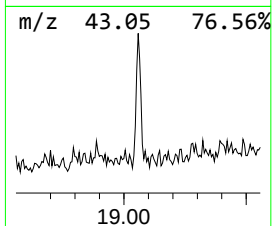
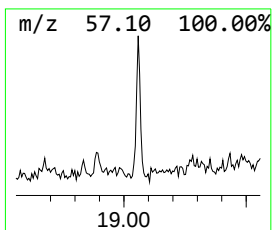
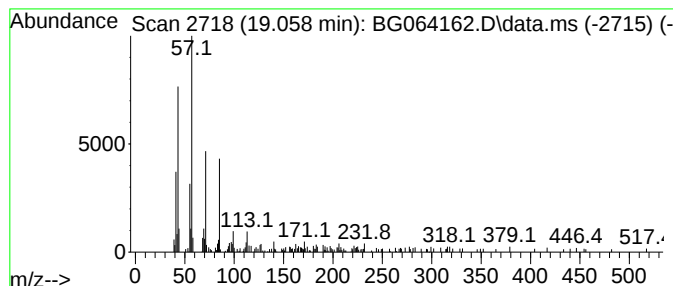
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.058	2.54 ng	80034	Phenanthrene-d10	17.225

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	91
2		Pentacosane	352	C25H52	000629-99-2	90
3		Heneicosane	296	C21H44	000629-94-7	78
4		Octadecane	254	C18H38	000593-45-3	72
5		Tetradecane, 2,5-dimethyl-	226	C16H34	056292-69-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040225\
Data File : BG064162.D
Acq On : 2 Apr 2025 21:04
Operator : RC/JU
Sample : Q1692-01 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
NB-08-04012025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG030525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.992	4.5	ng	61433	1	7.865	275461	20.0
Silane, trichlo...	12.084	2.9	ng	61110	2	10.650	419503	20.0
Tridecane	13.323	3.1	ng	84561	3	14.487	548995	20.0
Pentadecane	14.399	4.0	ng	109625	3	14.487	548995	20.0
Octadecane	15.351	4.0	ng	110327	3	14.487	548995	20.0
Heptadecane	16.208	4.1	ng	129613	4	17.225	631117	20.0
Nonadecane, 9-m...	17.002	3.4	ng	107668	4	17.225	631117	20.0
Nonadecane	17.736	3.2	ng	99948	4	17.225	631117	20.0
Pentacosane	18.429	2.8	ng	88215	4	17.225	631117	20.0
Heneicosane	19.058	2.5	ng	80034	4	17.225	631117	20.0