

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
 Data File : BG059761.D
 Acq On : 18 Nov 2023 12:31
 Operator : MA/JU
 Sample : 05401-01 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ETGI-338

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-BG111623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG059761.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.832	411	416	424	rVB2	68860	118424	17.39%	1.473%
2	7.236	651	655	661	rVB2	9053	9227	1.35%	0.115%
3	7.460	686	693	700	rBV	93908	168458	24.73%	2.095%
4	8.229	819	824	829	rVB3	10421	15490	2.27%	0.193%
5	8.341	835	843	850	rBV	143457	243882	35.81%	3.033%
6	8.494	865	869	872	rBV3	12591	18365	2.70%	0.228%
7	8.799	915	921	922	rBV2	15489	15818	2.32%	0.197%
8	8.928	938	943	951	rBV6	10559	22560	3.31%	0.281%
9	9.140	976	979	984	rBV2	6900	11765	1.73%	0.146%
10	9.539	1036	1047	1055	rVB3	60758	120570	17.70%	1.499%
11	11.190	1320	1328	1335	rBV2	199821	373510	54.84%	4.644%
12	12.495	1547	1550	1552	rBV3	12498	12018	1.76%	0.149%
13	12.571	1560	1563	1567	rVB3	6089	9491	1.39%	0.118%
14	13.047	1640	1644	1647	rBV5	12760	19984	2.93%	0.248%
15	13.082	1647	1650	1653	rVV4	6720	9874	1.45%	0.123%
16	13.118	1653	1656	1659	rVB4	10976	13840	2.03%	0.172%
17	13.218	1666	1673	1674	rBV3	10341	17706	2.60%	0.220%
18	13.294	1683	1686	1688	rBV3	11576	13354	1.96%	0.166%
19	13.347	1693	1695	1697	rBV3	9692	9055	1.33%	0.113%
20	13.423	1703	1708	1714	rVB9	16142	31780	4.67%	0.395%
21	13.529	1724	1726	1731	rBV6	9966	16105	2.36%	0.200%
22	13.594	1731	1737	1744	rBV2	243639	394255	57.88%	4.902%
23	13.717	1753	1758	1763	rBV3	57559	87106	12.79%	1.083%
24	13.805	1767	1773	1775	rBV6	13170	23261	3.42%	0.289%
25	13.840	1777	1779	1781	rBV3	10314	11724	1.72%	0.146%
26	13.893	1786	1788	1789	rVV2	11977	12190	1.79%	0.152%
27	13.917	1791	1792	1795	rBV3	10267	9405	1.38%	0.117%
28	13.940	1795	1796	1798	rBV2	12357	10219	1.50%	0.127%
29	14.005	1802	1807	1808	rBV4	9451	13013	1.91%	0.162%
30	14.064	1814	1817	1819	rBV4	12768	14858	2.18%	0.185%
31	14.281	1850	1854	1857	rBV6	14785	20282	2.98%	0.252%
32	14.357	1863	1867	1872	rVB6	40350	65206	9.57%	0.811%
33	14.404	1872	1875	1878	rBV5	16264	20520	3.01%	0.255%
34	14.451	1880	1883	1886	rBV5	11722	15596	2.29%	0.194%
35	14.522	1891	1895	1898	rVV5	19969	23980	3.52%	0.298%
36	14.745	1931	1933	1934	rBV2	13208	9385	1.38%	0.117%

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Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.780	1934	1939	1942	rVV3	111602	156440	22.97%	1.945%
38	14.857	1950	1952	1955	rVB4	17388	17373	2.55%	0.216%
39	14.904	1957	1960	1962	rVV4	15970	21423	3.15%	0.266%
40	14.968	1966	1971	1978	rVB	411620	653864	96.00%	8.131%
41	15.151	1999	2002	2007	rVB7	13726	20361	2.99%	0.253%
42	15.274	2020	2023	2025	rBV4	12293	13927	2.04%	0.173%
43	15.339	2030	2034	2038	rVB7	26609	29803	4.38%	0.371%
44	15.386	2040	2042	2045	rVB4	26807	23517	3.45%	0.292%
45	15.421	2045	2048	2050	rBV3	25699	28724	4.22%	0.357%
46	15.550	2067	2070	2071	rBV3	22880	19882	2.92%	0.247%
47	15.726	2095	2100	2104	rVB3	134853	167225	24.55%	2.079%
48	15.785	2104	2110	2114	rBV9	25117	54287	7.97%	0.675%
49	15.850	2116	2121	2122	rBV5	19199	20236	2.97%	0.252%
50	15.920	2130	2133	2138	rBV6	19626	36007	5.29%	0.448%
51	16.014	2146	2149	2150	rBV3	10631	11682	1.72%	0.145%
52	16.126	2161	2168	2175	rBV3	61160	128156	18.82%	1.594%
53	16.220	2182	2184	2186	rVB3	12507	10139	1.49%	0.126%
54	16.284	2193	2195	2196	rVB2	16935	11019	1.62%	0.137%
55	16.349	2204	2206	2209	rBV4	22667	32683	4.80%	0.406%
56	16.455	2218	2224	2229	rBV3	184745	313645	46.05%	3.900%
57	16.590	2242	2247	2254	rBV2	196117	343186	50.39%	4.267%
58	16.866	2292	2294	2295	rVB2	19628	11178	1.64%	0.139%
59	16.925	2301	2304	2307	rVB5	22797	35384	5.20%	0.440%
60	16.954	2307	2309	2310	rBV2	16477	11104	1.63%	0.138%
61	16.984	2312	2314	2318	rVB5	25289	28143	4.13%	0.350%
62	17.048	2322	2325	2327	rBV4	17485	19066	2.80%	0.237%
63	17.136	2338	2340	2342	rBV3	13386	11929	1.75%	0.148%
64	17.160	2342	2344	2349	rVB6	14977	17922	2.63%	0.223%
65	17.213	2351	2353	2354	rBV2	14665	9045	1.33%	0.112%
66	17.383	2378	2382	2386	rVV2	148958	187696	27.56%	2.334%
67	17.430	2386	2390	2394	rVB6	49863	78306	11.50%	0.974%
68	17.501	2400	2402	2405	rBV4	10875	13258	1.95%	0.165%
69	17.677	2430	2432	2433	rBV2	16461	15389	2.26%	0.191%
70	17.724	2434	2440	2445	rVV	435618	681107	100.00%	8.469%
71	17.836	2457	2459	2460	rBV2	16212	13151	1.93%	0.164%
72	17.865	2460	2464	2466	rBV5	16711	24860	3.65%	0.309%
73	18.124	2503	2508	2513	rVV2	142550	183186	26.90%	2.278%
74	18.171	2515	2516	2518	rVB2	19811	13287	1.95%	0.165%
75	18.235	2523	2527	2529	rBV4	19121	21809	3.20%	0.271%
76	18.312	2538	2540	2545	rVB6	51953	57634	8.46%	0.717%

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77	18.400	2552	2555	2556	rBV3	24725	23106	3.39%	0.287%
78	18.529	2571	2577	2582	rBV9	52727	99372	14.59%	1.236%
79	18.629	2592	2594	2595	rBV2	19533	13298	1.95%	0.165%
80	18.811	2621	2625	2629	rVB	117380	129554	19.02%	1.611%
81	19.134	2678	2680	2681	rBV2	17486	13685	2.01%	0.170%
82	19.440	2727	2732	2735	rBV5	159441	220097	32.31%	2.737%
83	19.598	2756	2759	2763	rBV6	22837	35904	5.27%	0.446%
84	19.751	2783	2785	2787	rBV3	21397	15640	2.30%	0.194%
85	19.792	2790	2792	2794	rVV3	24396	16276	2.39%	0.202%
86	20.004	2825	2828	2832	rBV5	61939	69574	10.21%	0.865%
87	20.292	2873	2877	2884	rBV	344067	480182	70.50%	5.971%
88	20.515	2913	2915	2916	rBV2	29095	24126	3.54%	0.300%
89	21.854	3140	3143	3148	rVB7	42748	57138	8.39%	0.710%
90	21.966	3161	3162	3164	rBV2	31430	18211	2.67%	0.226%
91	22.048	3171	3176	3182	rBV2	334987	576033	84.57%	7.163%
92	25.644	3781	3788	3799	rVB2	184569	566174	83.13%	7.040%
93	26.943	4007	4009	4010	rBV	27815	18086	2.66%	0.225%
94	30.368	4591	4592	4593	rVB	32618	11498	1.69%	0.143%
95	30.973	4693	4695	4697	rBV3	21965	19425	2.85%	0.242%
96	31.114	4717	4719	4722	rBV4	23535	31536	4.63%	0.392%
97	31.578	4797	4798	4800	rBV2	25390	11384	1.67%	0.142%
98	31.619	4803	4805	4807	rBV3	17829	16156	2.37%	0.201%
99	31.995	4867	4869	4872	rBV4	20492	14877	2.18%	0.185%
100	32.366	4930	4932	4934	rBV3	22178	11398	1.67%	0.142%

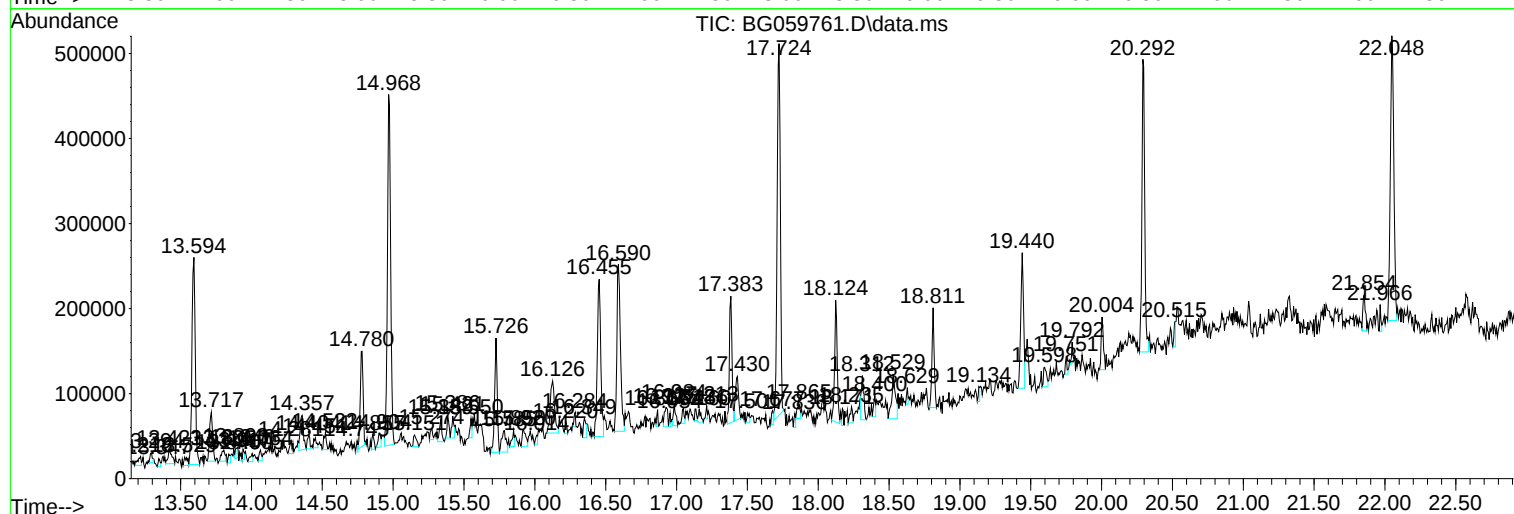
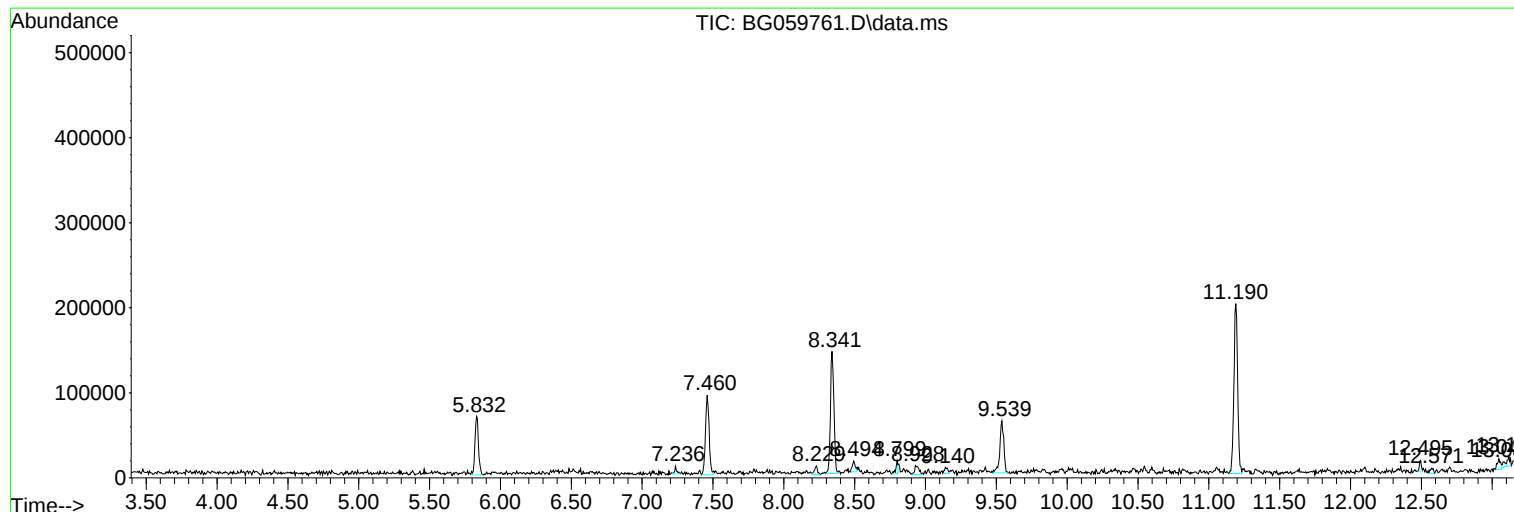
Sum of corrected areas: 8042039

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Instrument :
BNA_G
ClientSampleId :
ETGI-338

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.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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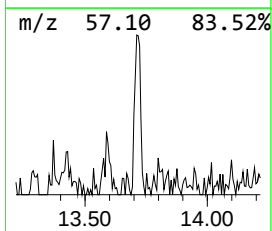
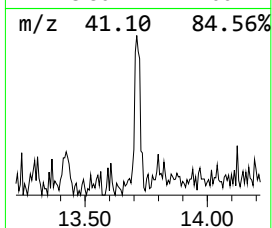
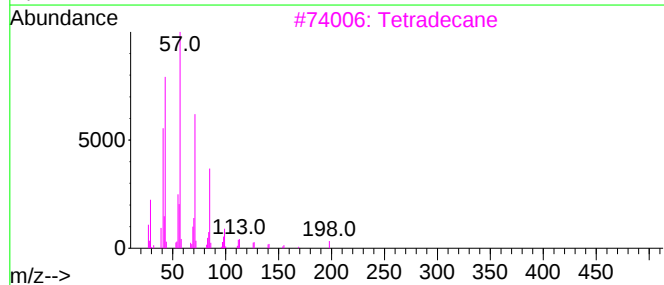
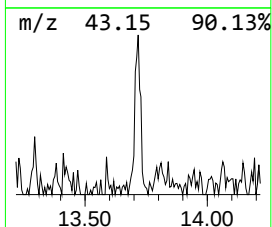
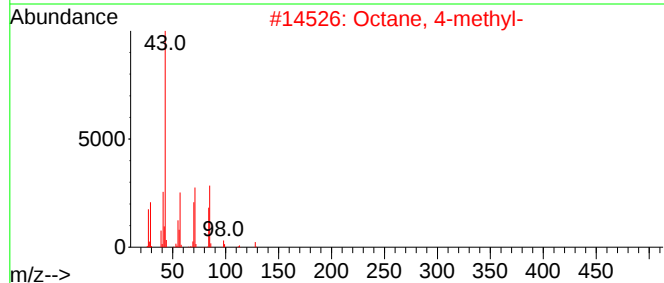
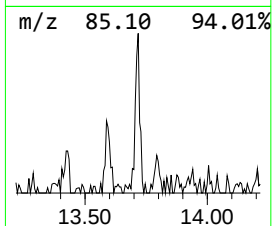
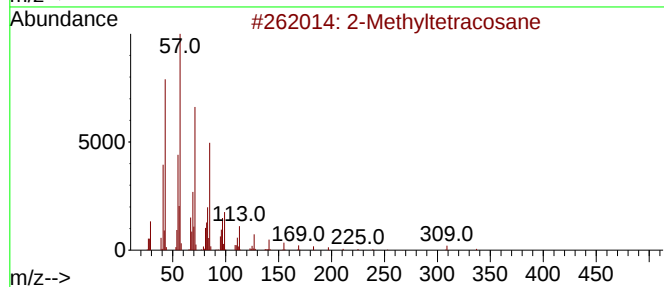
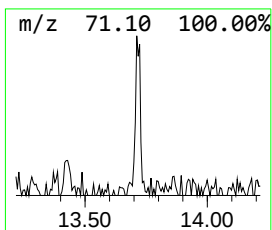
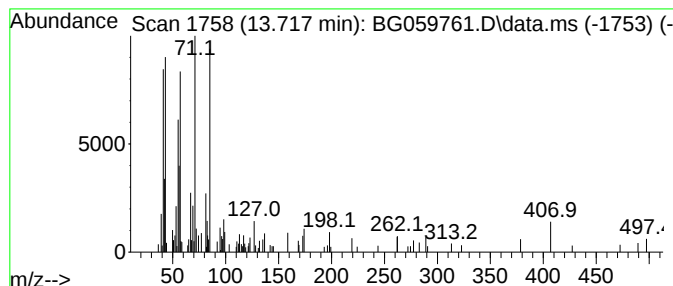
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Methyltetracosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.717	2.66 ng	87106	Acenaphthene-d10	14.968

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Methyltetracosane	352	C25H52	001560-78-7	64
2		Octane, 4-methyl-	128	C9H20	002216-34-4	64
3		Tetradecane	198	C14H30	000629-59-4	50
4		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50
5		2,6-Octadiene-4,5-diol, 4-methyl-	156	C9H16O2	056335-74-1	50



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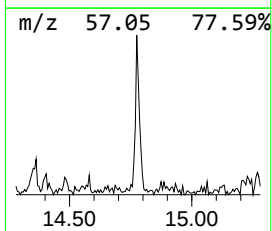
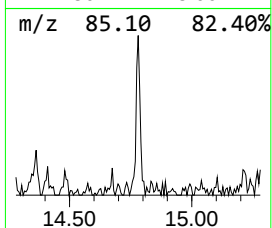
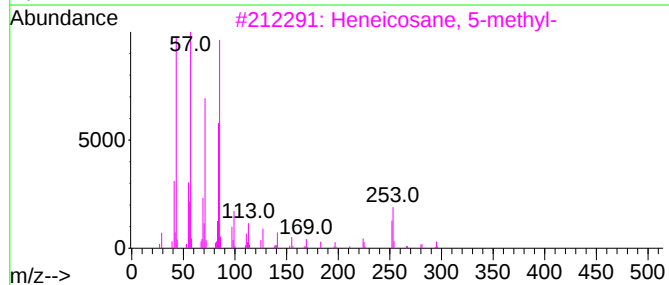
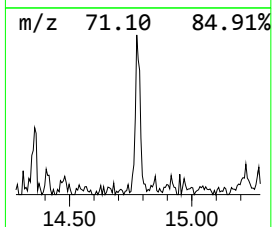
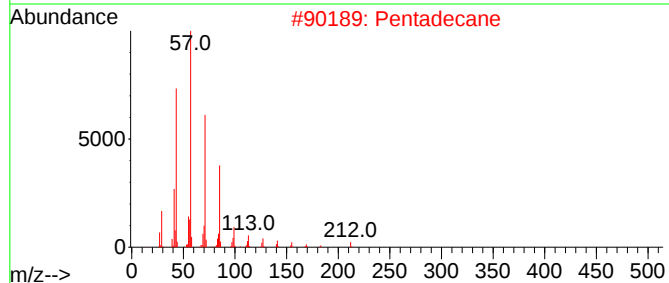
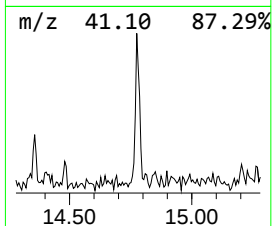
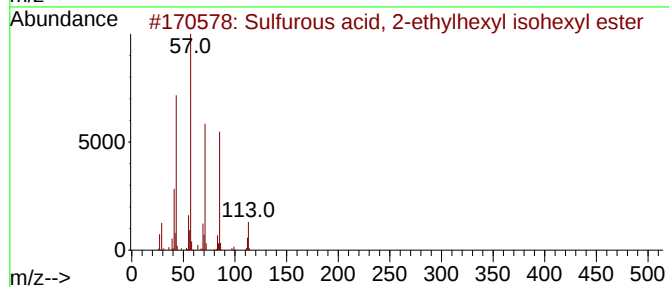
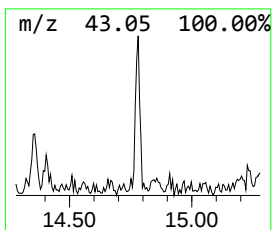
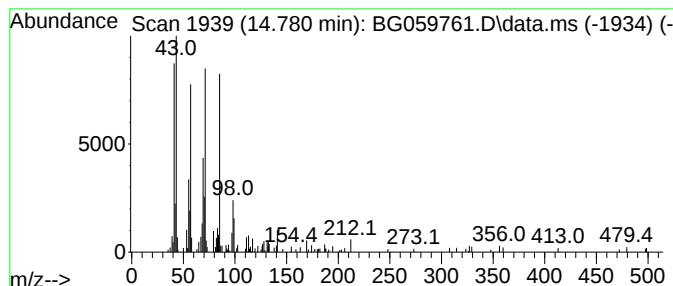
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Sulfurous acid, 2-ethylhexy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	4.79 ng	156440	Acenaphthene-d10	14.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	59
2			Pentadecane	212	C15H32	000629-62-9	52
3			Heneicosane, 5-methyl-	310	C22H46	025117-37-7	47
4			2-Bromononane	206	C9H19Br	002216-35-5	43
5			Eicosane, 2,4-dimethyl-	310	C22H46	075163-98-3	38



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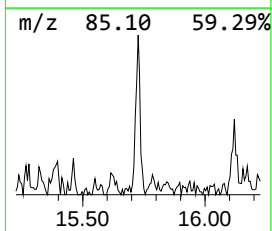
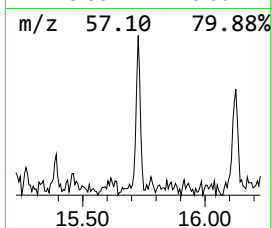
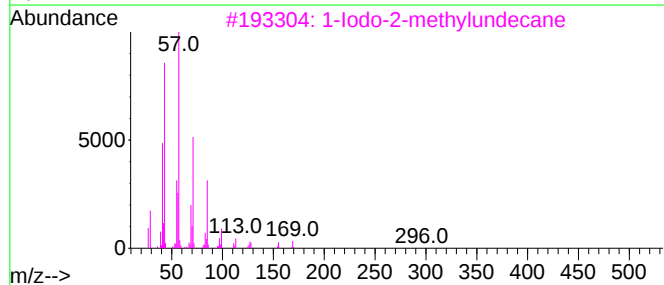
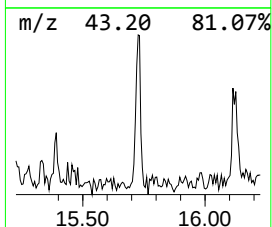
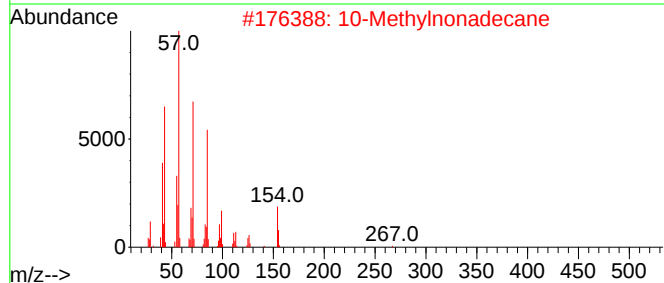
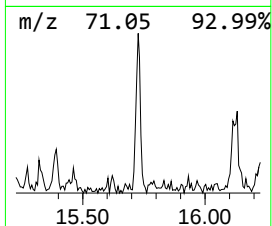
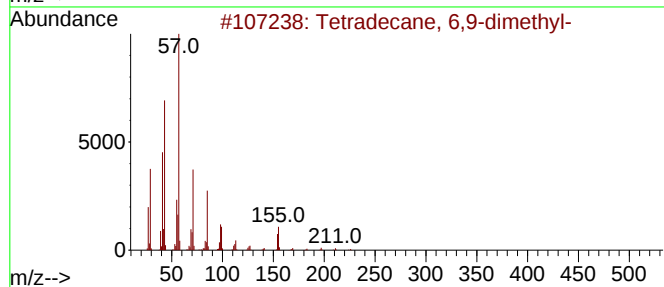
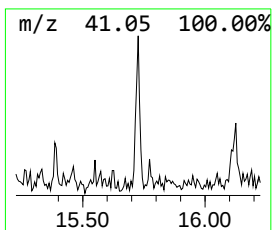
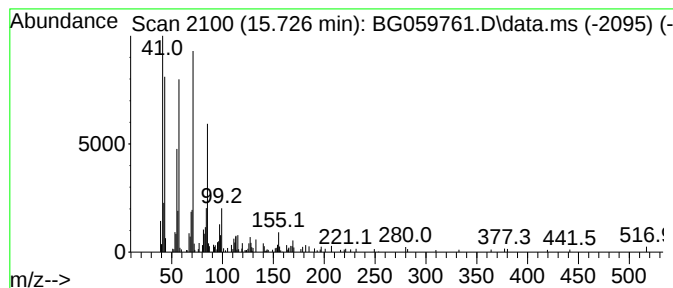
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tetradecane, 6,9-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.726	5.11 ng	167225	Acenaphthene-d10	14.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 6,9-dimethyl-	226	C16H34	055045-13-1	81
2			10-Methylnonadecane	282	C20H42	056862-62-5	80
3			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	80
4			Pentacosane	352	C25H52	000629-99-2	80
5			Pentadecane, 4-methyl-	226	C16H34	002801-87-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
Operator : MA/JU
Sample : 05401-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-338

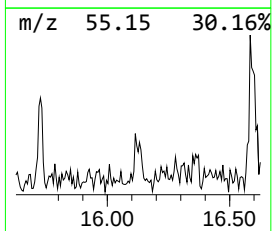
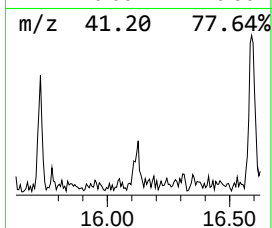
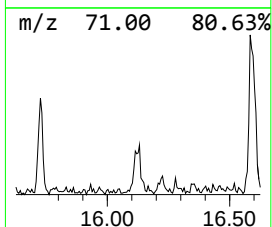
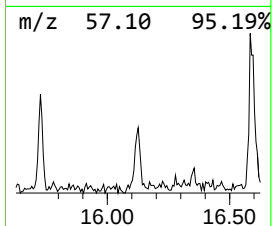
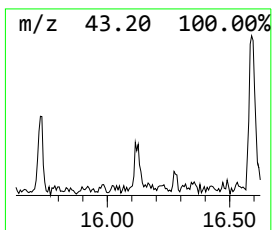
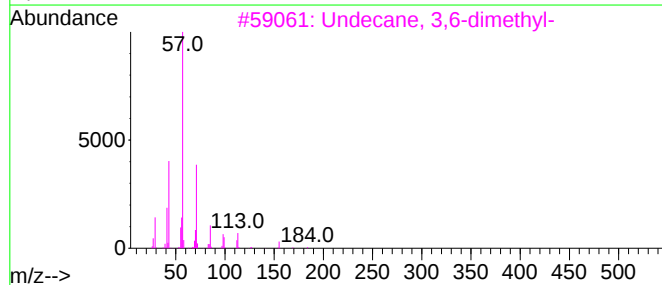
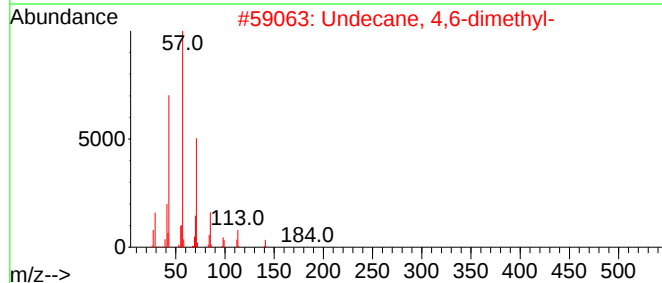
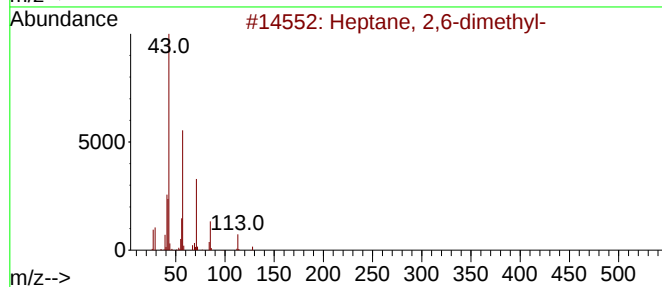
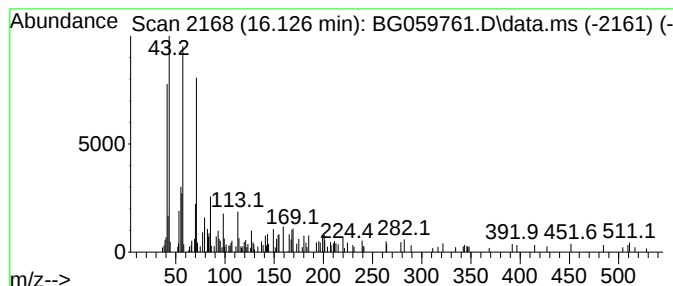
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Heptane, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.126	3.92 ng	128156	Acenaphthene-d10	14.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	76
2			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	64
3			Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	64
4			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	59
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
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ALS Vial : 26 Sample Multiplier: 1

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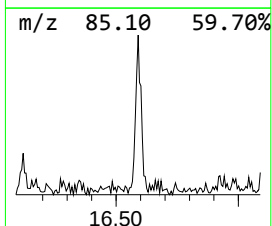
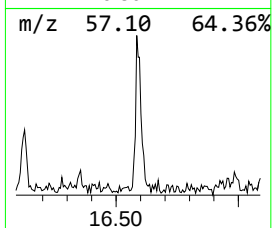
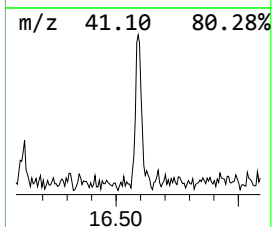
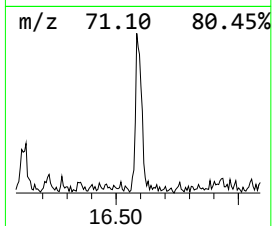
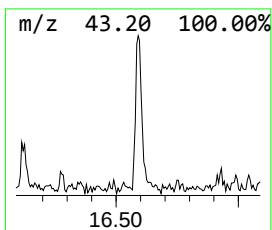
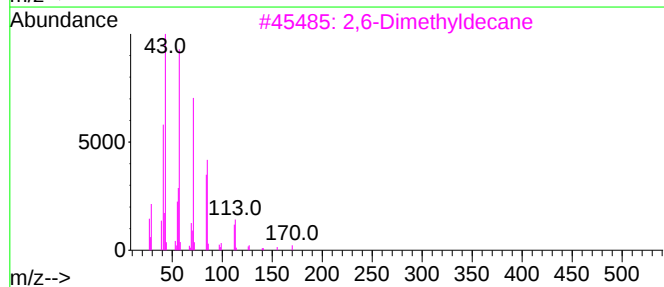
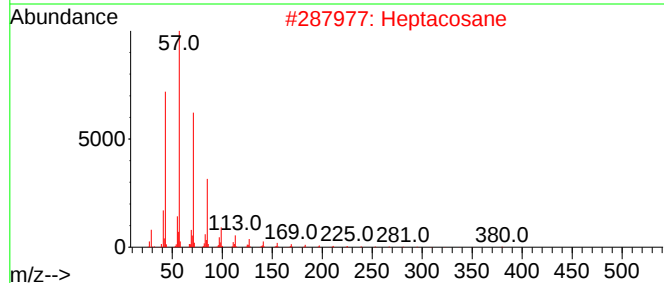
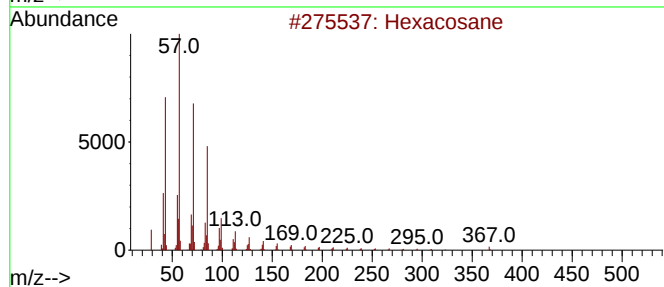
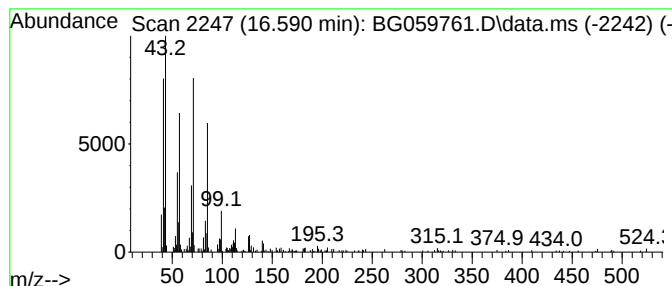
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexacosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.590	10.08 ng	343186	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	80
2			Heptacosane	380	C27H56	000593-49-7	59
3			2,6-Dimethyldecane	170	C12H26	013150-81-7	58
4			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	58
5			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
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Misc :
ALS Vial : 26 Sample Multiplier: 1

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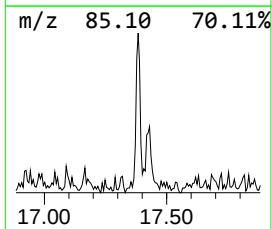
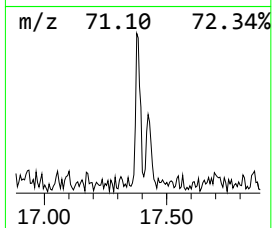
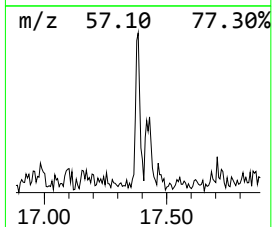
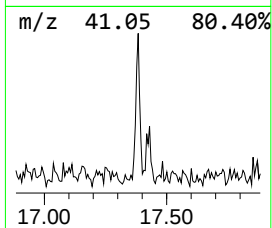
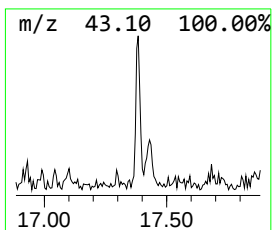
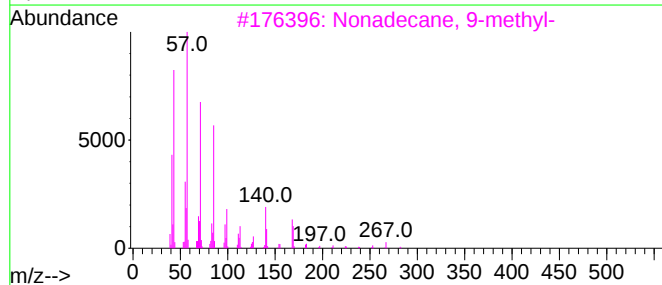
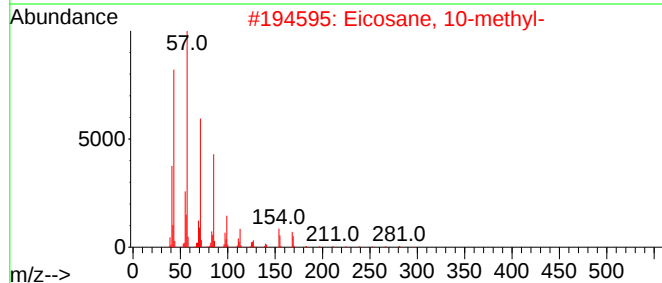
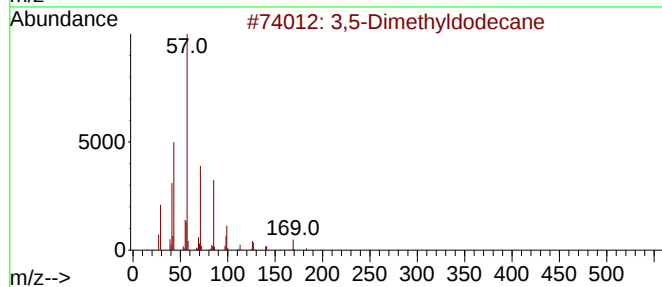
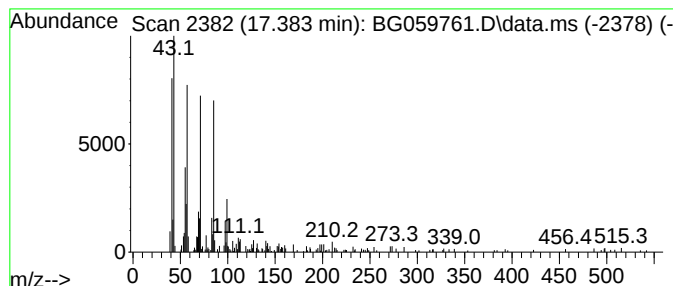
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 3,5-Dimethyldodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.383	5.51 ng	187696	Phenanthrene-d10	17.718

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,5-Dimethyldodecane	198	C14H30	107770-99-0	80
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
3		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	80
4		Octacosane	394	C28H58	000630-02-4	64
5		Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
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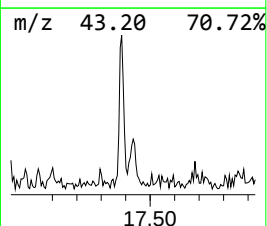
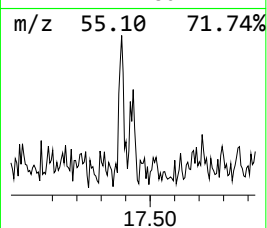
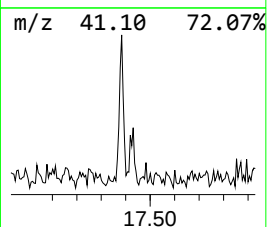
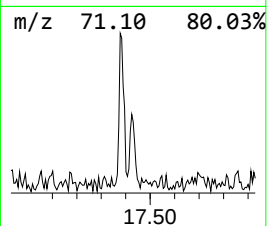
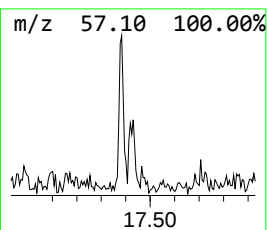
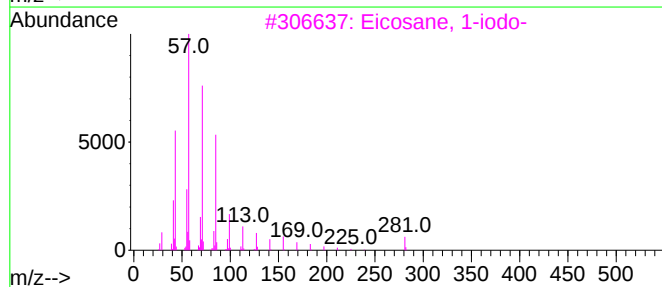
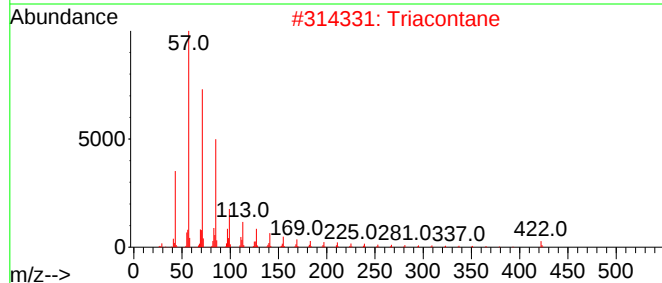
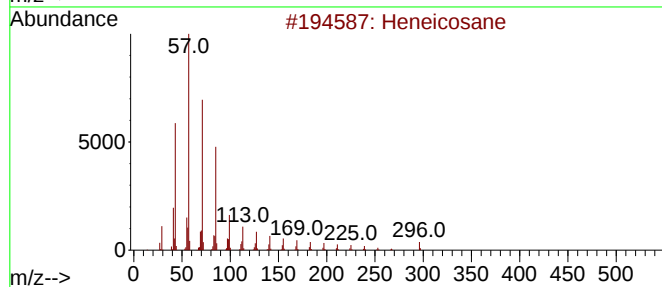
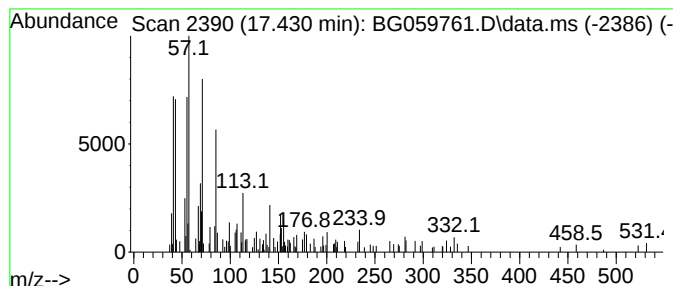
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.430	2.30 ng	78306	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	52
2			Triacontane	422	C30H62	000638-68-6	50
3			Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	50
4			Nonadecane	268	C19H40	000629-92-5	49
5			Heptadecane	240	C17H36	000629-78-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
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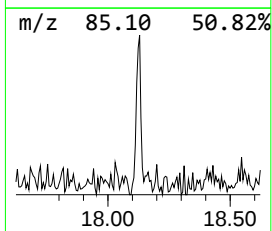
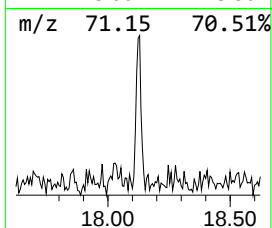
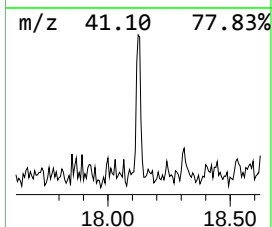
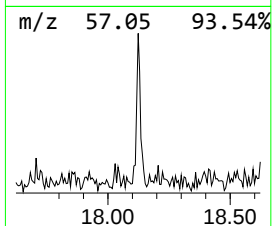
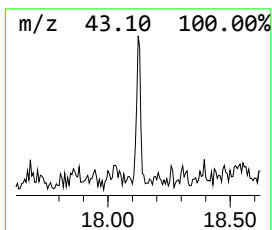
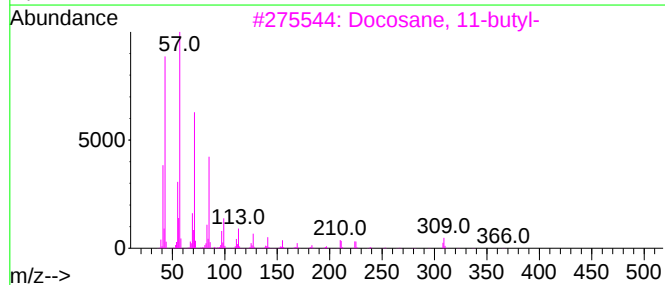
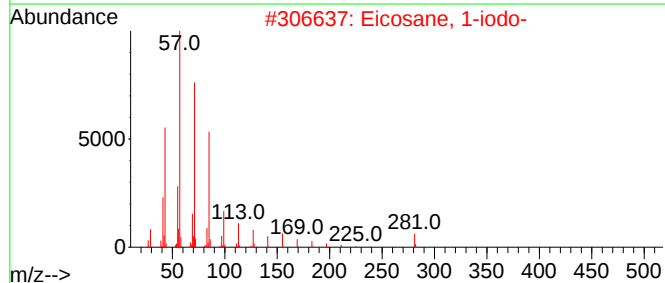
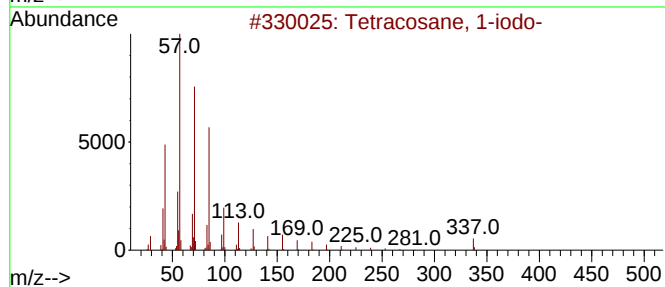
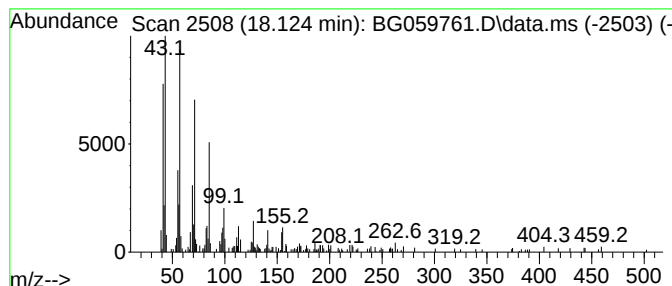
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetracosane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.124	5.38 ng	183186	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	74
2			Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	72
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	72
4			Heneicosane	296	C21H44	000629-94-7	72
5			2-Bromotetradecane	276	C14H29Br	074036-95-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
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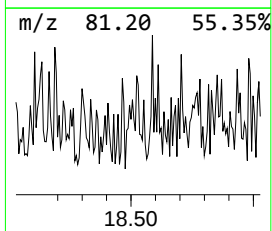
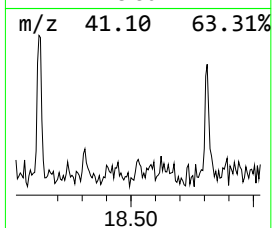
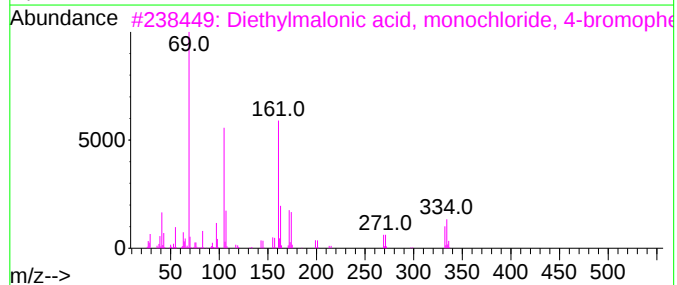
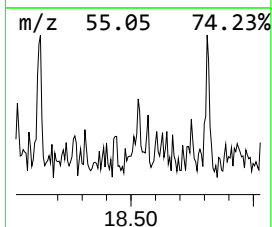
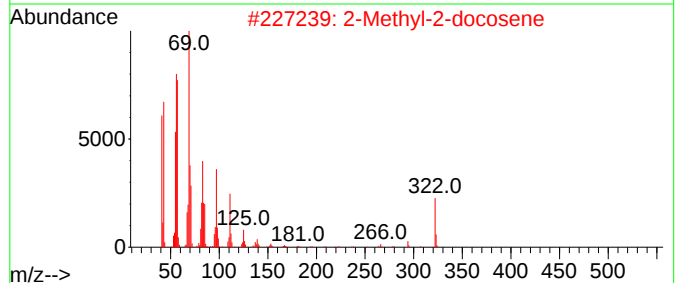
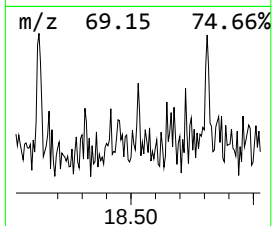
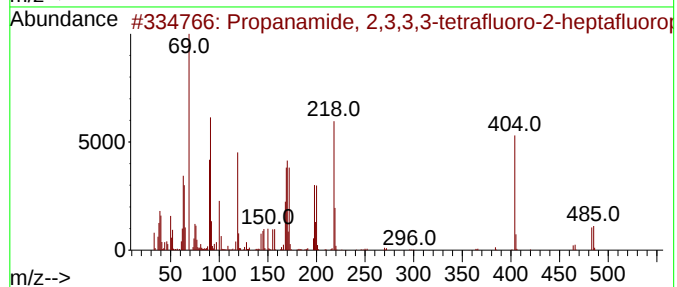
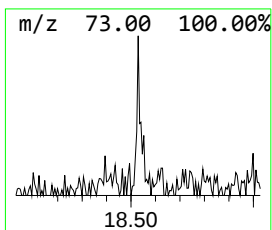
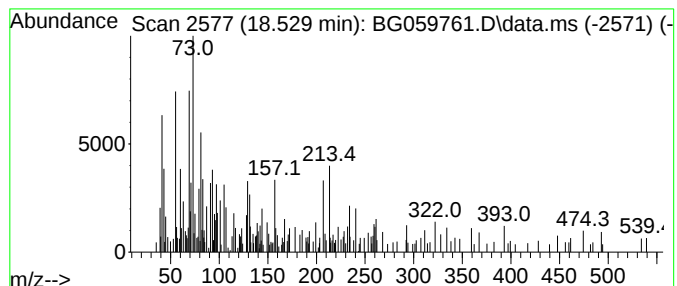
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown18.529 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.529	2.92 ng	99372	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanamide, 2,3,3,3-tetrafluoro...	483	C12H5BrF11N02	294195-06-5	22
2			2-Methyl-2-docosene	322	C23H46	1000131-16-9	22
3			Diethylmalonic acid, monochlorid...	332	C13H14BrCl03	1000369-83-6	22
4			Hexafluoroisopropyl acrylate	222	C6H4F6O2	002160-89-6	22
5			4,25-Secoobscurinervan-4-ol, 6,7...	428	C24H32N2O5	007236-84-2	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
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ClientSampleId :
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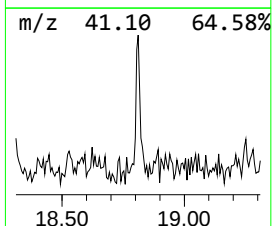
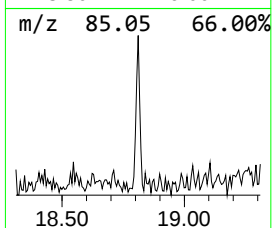
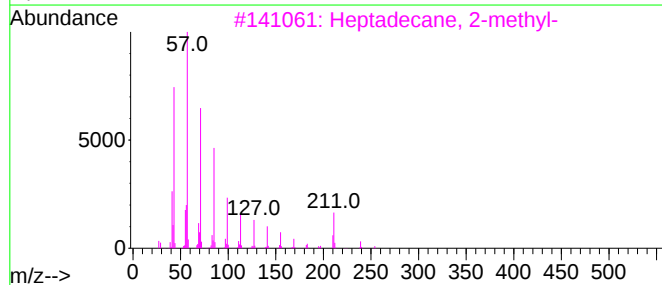
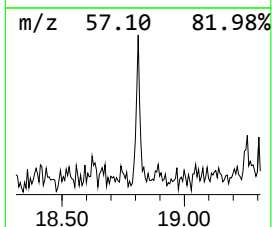
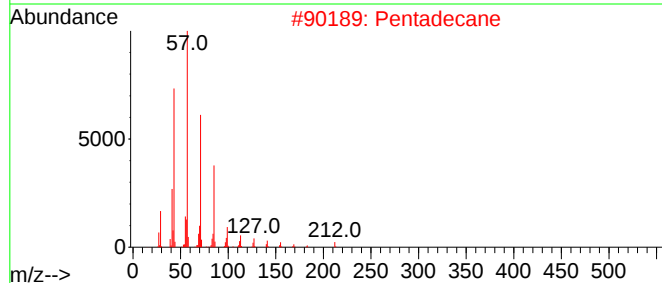
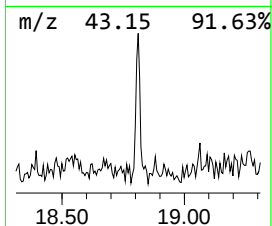
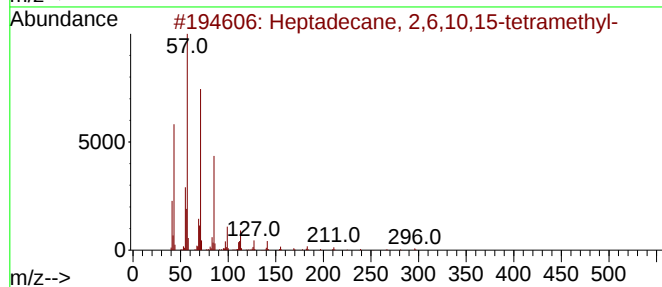
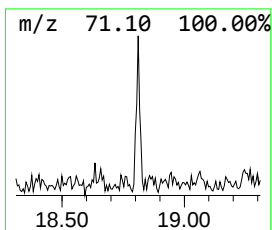
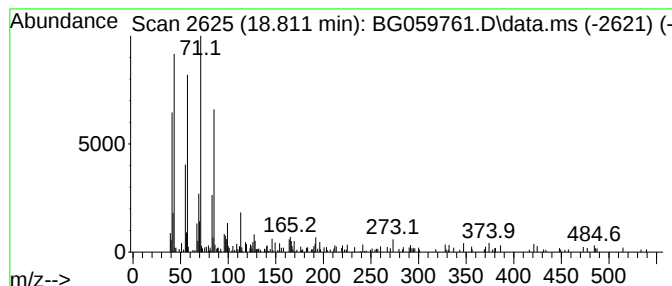
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 2,6,10,15-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.811	3.80 ng	129554	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
2			Pentadecane	212	C15H32	000629-62-9	83
3			Heptadecane, 2-methyl-	254	C18H38	001560-89-0	80
4			Nonadecane	268	C19H40	000629-92-5	80
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
Operator : MA/JU
Sample : 05401-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-338

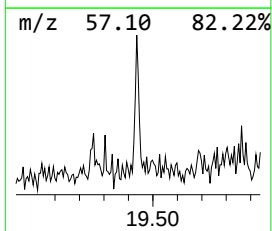
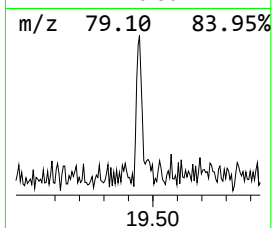
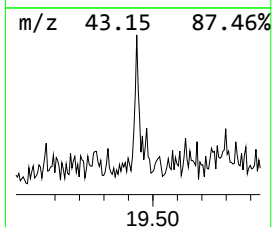
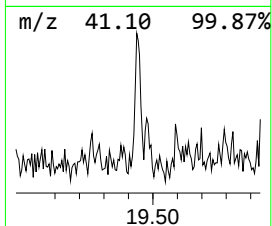
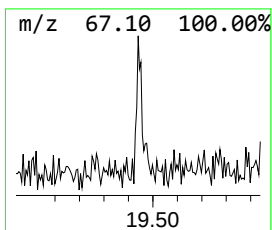
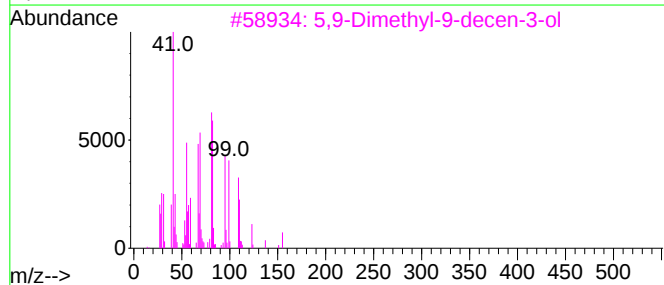
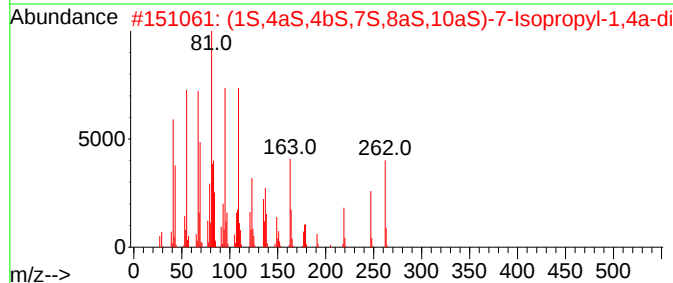
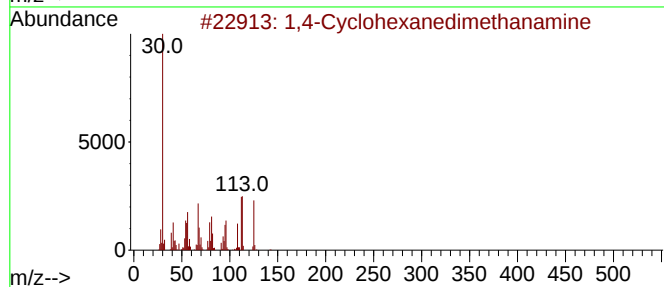
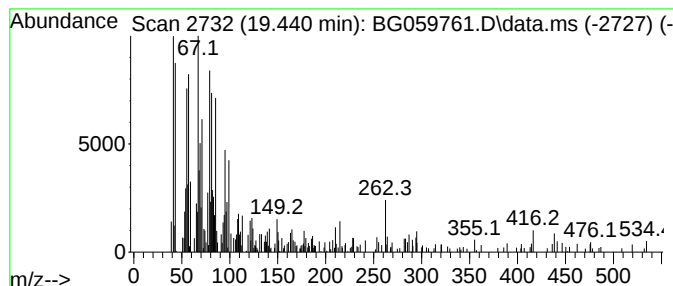
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown19.440 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.440	6.46 ng	220097	Phenanthrene-d10	17.718

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Cyclohexanedimethanamine	142	C8H18N2	002549-93-1	47
2			(1S,4aS,4bS,7S,8aS,10aS)-7-Isopr...	262	C19H34	002221-95-6	41
3			5,9-Dimethyl-9-decen-3-ol	184	C12H24O	118447-56-6	38
4			8-Decen-3-ol, 5,9-dimethyl-	184	C12H24O	019550-54-0	38
5			6-(2-propynyloxy)-1-hexanol	156	C9H16O2	058757-86-1	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
Operator : MA/JU
Sample : 05401-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ETGI-338

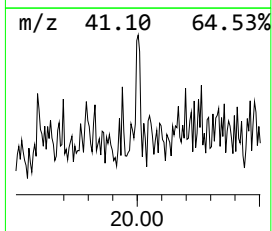
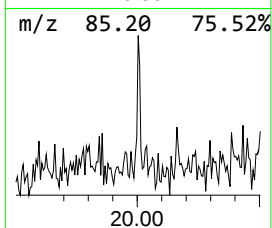
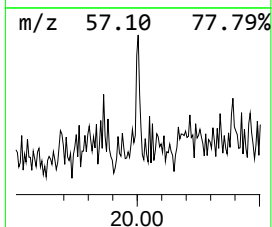
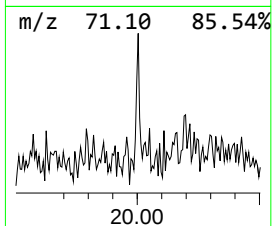
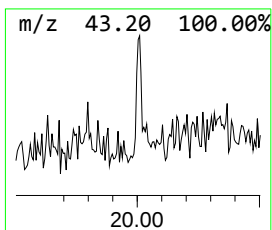
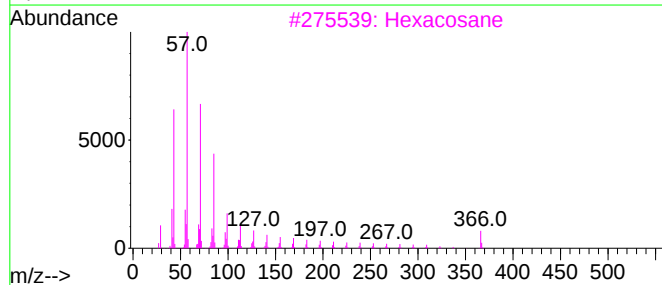
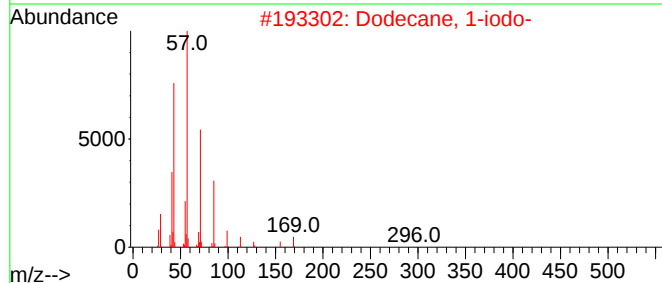
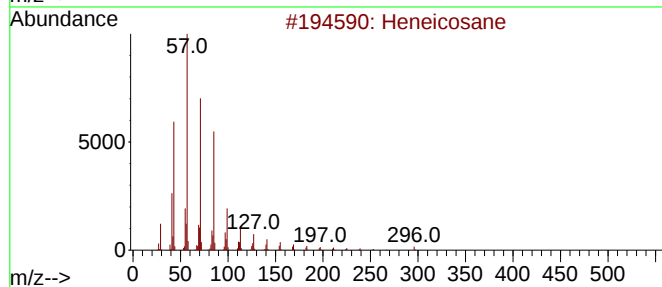
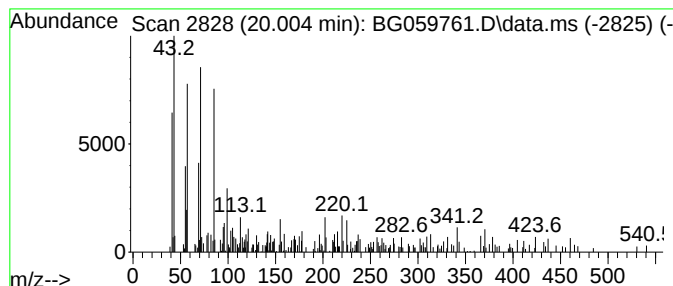
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Dodecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.004	2.42 ng	69574	Chrysene-d12	22.048

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	74
2		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
3		Hexacosane	366	C26H54	000630-01-3	72
4		Octacosane	394	C28H58	000630-02-4	72
5		Pentadecane	212	C15H32	000629-62-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
Data File : BG059761.D
Acq On : 18 Nov 2023 12:31
Operator : MA/JU
Sample : 05401-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ETGI-338

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.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-Methyltetraco...	13.717	2.7	ng	87106	3	14.968	653864	20.0
Sulfurous acid,...	14.780	4.8	ng	156440	3	14.968	653864	20.0
Tetradecane, 6,...	15.726	5.1	ng	167225	3	14.968	653864	20.0
Heptane, 2,6-di...	16.126	3.9	ng	128156	3	14.968	653864	20.0
Hexacosane	16.590	10.1	ng	343186	4	17.718	681107	20.0
3,5-Dimethyldod...	17.383	5.5	ng	187696	4	17.718	681107	20.0
Heneicosane	17.430	2.3	ng	78306	4	17.718	681107	20.0
Tetracosane, 1-...	18.124	5.4	ng	183186	4	17.718	681107	20.0
unknown18.529	18.529	2.9	ng	99372	4	17.718	681107	20.0
Heptadecane, 2,...	18.811	3.8	ng	129554	4	17.718	681107	20.0
unknown19.440	19.440	6.5	ng	220097	4	17.718	681107	20.0
Dodecane, 1-iodo-	20.004	2.4	ng	69574	5	22.048	576033	20.0