

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
 Data File : BG063108.D
 Acq On : 28 Sep 2024 15:06
 Operator : RC/JU
 Sample : P3927-13DL 30X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DD00DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\SFAM-EPA-BG092524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.875	470	474	477	rBV3	11693	13596	0.61%	0.117%
2	6.422	562	567	568	rBV2	5835	6328	0.28%	0.055%
3	6.839	630	638	643	rBV4	11628	24120	1.08%	0.208%
4	6.898	646	648	652	rBV4	7354	9984	0.45%	0.086%
5	6.939	652	655	661	rVB	16110	25399	1.14%	0.219%
6	7.003	664	666	674	rVB7	16381	28942	1.30%	0.250%
7	7.080	676	679	686	rVB3	9502	13053	0.58%	0.113%
8	7.186	692	697	703	rBV7	10061	15981	0.72%	0.138%
9	7.244	703	707	709	rBV2	9953	13889	0.62%	0.120%
10	7.268	709	711	712	rBV2	9885	6128	0.27%	0.053%
11	7.380	727	730	731	rBV2	6889	6553	0.29%	0.057%
12	7.479	742	747	757	rVB5	37198	97669	4.37%	0.844%
13	7.556	757	760	762	rBV2	7993	7447	0.33%	0.064%
14	7.850	803	810	816	rBV	226896	389050	17.41%	3.362%
15	7.955	827	828	830	rBV2	6956	5929	0.27%	0.051%
16	8.114	851	855	858	rVV3	5669	7628	0.34%	0.066%
17	8.143	858	860	863	rVB4	6967	6455	0.29%	0.056%
18	8.273	879	882	885	rVB5	9889	10701	0.48%	0.092%
19	8.331	885	892	894	rBV6	10222	16005	0.72%	0.138%
20	8.384	898	901	903	rBV3	8975	10814	0.48%	0.093%
21	8.490	914	919	920	rBV4	16276	18684	0.84%	0.161%
22	8.560	928	931	936	rVB3	10043	13342	0.60%	0.115%
23	8.613	936	940	941	rBV2	12175	9637	0.43%	0.083%
24	8.796	968	971	972	rVV2	11136	9539	0.43%	0.082%
25	8.848	975	980	984	rVB7	16182	28637	1.28%	0.247%
26	8.931	990	994	995	rBV3	5678	7798	0.35%	0.067%
27	8.960	995	999	1002	rVB4	13868	19816	0.89%	0.171%
28	9.078	1013	1019	1024	rVB5	32055	52386	2.34%	0.453%
29	9.636	1112	1114	1121	rVB5	7360	6726	0.30%	0.058%
30	9.736	1126	1131	1135	rBV6	8271	17273	0.77%	0.149%
31	9.777	1136	1138	1142	rVB4	5944	7068	0.32%	0.061%
32	10.006	1172	1177	1180	rVV5	7334	14798	0.66%	0.128%
33	10.106	1191	1194	1196	rVB2	4771	5945	0.27%	0.051%
34	10.276	1220	1223	1226	rVV3	10752	13890	0.62%	0.120%
35	10.329	1231	1232	1237	rVB4	8650	11955	0.54%	0.103%
36	10.635	1272	1284	1291	rBV	390870	746519	33.41%	6.451%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\SFAM-EPA-BG092524.MA.M
 Title : SVOA CALIBRATION

37	10.758	1300	1305	1311	rBV9	24537	53355	2.39%	0.461%
38	11.016	1343	1349	1355	rVB3	17272	34552	1.55%	0.299%
39	11.216	1379	1383	1386	rVB	7553	9310	0.42%	0.080%
40	11.528	1432	1436	1437	rBV3	7876	11342	0.51%	0.098%
41	11.669	1455	1460	1464	rBV2	7797	11721	0.52%	0.101%
42	11.880	1494	1496	1498	rVV3	14641	13219	0.59%	0.114%
43	11.904	1498	1500	1505	rVB4	11135	13444	0.60%	0.116%
44	12.056	1521	1526	1528	rBV3	13636	19338	0.87%	0.167%
45	12.098	1531	1533	1538	rVB3	8146	10190	0.46%	0.088%
46	12.315	1565	1570	1575	rVB5	14854	23602	1.06%	0.204%
47	12.479	1591	1598	1602	rBV5	16084	37551	1.68%	0.324%
48	12.515	1602	1604	1608	rVV3	7189	9054	0.41%	0.078%
49	12.996	1683	1686	1687	rBV	5108	5908	0.26%	0.051%
50	13.055	1689	1696	1699	rVV5	25588	42553	1.90%	0.368%
51	13.308	1735	1739	1744	rBV4	8502	13582	0.61%	0.117%
52	13.525	1773	1776	1779	rVV3	9813	13777	0.62%	0.119%
53	13.666	1796	1800	1805	rVV7	14373	24588	1.10%	0.212%
54	13.813	1819	1825	1826	rBV3	10155	14948	0.67%	0.129%
55	13.854	1830	1832	1833	rVV2	6845	6509	0.29%	0.056%
56	13.890	1834	1838	1843	rVV	22103	36935	1.65%	0.319%
57	14.036	1858	1863	1867	rBV7	9264	15423	0.69%	0.133%
58	14.177	1883	1887	1891	rVV2	26237	41144	1.84%	0.356%
59	14.207	1891	1892	1899	rVB3	12054	13139	0.59%	0.114%
60	14.389	1918	1923	1928	rBV5	17680	27241	1.22%	0.235%
61	14.483	1933	1939	1949	rVB	691008	1082144	48.43%	9.351%
62	14.630	1960	1964	1966	rVB4	4915	6255	0.28%	0.054%
63	14.730	1980	1981	1986	rVB3	9189	9786	0.44%	0.085%
64	15.065	2035	2038	2041	rBV4	5374	6598	0.30%	0.057%
65	15.118	2041	2047	2053	rBV5	56736	103074	4.61%	0.891%
66	15.276	2072	2074	2075	rVV2	7674	5983	0.27%	0.052%
67	15.311	2078	2080	2083	rBV4	6939	7734	0.35%	0.067%
68	15.476	2104	2108	2114	rBV2	44091	70720	3.16%	0.611%
69	15.629	2130	2134	2140	rBV4	17245	34894	1.56%	0.302%
70	15.746	2152	2154	2159	rVB4	6446	9990	0.45%	0.086%
71	15.846	2167	2171	2178	rVB5	18273	34456	1.54%	0.298%
72	16.199	2229	2231	2235	rBV4	13286	18340	0.82%	0.158%
73	16.545	2287	2290	2294	rBV6	6708	10750	0.48%	0.093%
74	16.604	2296	2300	2301	rBV4	9744	9251	0.41%	0.080%
75	16.733	2319	2322	2325	rVV5	6859	9429	0.42%	0.081%
76	16.892	2342	2349	2361	rBV	664257	1186406	53.10%	10.252%

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 ClientSampleId :
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Integration Parameters: LSCINT.P
 Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BG092524.MA.M
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77	17.233	2400	2407	2417	rBV	1106014	1647262	73.72%	14.235%
78	17.333	2419	2424	2428	rVV2	53270	83156	3.72%	0.719%
79	17.485	2446	2450	2453	rVB4	9010	10937	0.49%	0.095%
80	17.526	2453	2457	2464	rBV8	8513	22240	1.00%	0.192%
81	17.744	2490	2494	2498	rVB5	12568	18255	0.82%	0.158%
82	17.808	2498	2505	2506	rBV7	10384	21670	0.97%	0.187%
83	17.855	2511	2513	2514	rBV2	10799	6193	0.28%	0.054%
84	17.908	2519	2522	2526	rVV4	19868	32578	1.46%	0.282%
85	17.967	2529	2532	2533	rVV3	7478	6849	0.31%	0.059%
86	18.378	2599	2602	2606	rBV5	8564	10578	0.47%	0.091%
87	18.596	2638	2639	2644	rVB5	10573	10297	0.46%	0.089%
88	19.066	2717	2719	2721	rBV3	10612	12148	0.54%	0.105%
89	19.224	2742	2746	2748	rBV4	14290	16058	0.72%	0.139%
90	19.283	2753	2756	2760	rBV4	5913	8153	0.36%	0.070%
91	19.318	2760	2762	2764	rBV3	8857	6129	0.27%	0.053%
92	19.542	2798	2800	2806	rVB7	15048	19370	0.87%	0.167%
93	19.630	2811	2815	2820	rVV	65943	99926	4.47%	0.864%
94	20.235	2916	2918	2924	rVB7	25710	31531	1.41%	0.272%
95	20.870	3022	3026	3031	rVB5	43026	62887	2.81%	0.543%
96	21.486	3125	3131	3143	rVB2	1358460	2145430	96.02%	18.540%
97	22.227	3252	3257	3261	rBV7	71194	125661	5.62%	1.086%
98	23.038	3393	3395	3402	rVB6	66127	119350	5.34%	1.031%
99	24.530	3640	3649	3664	rVB	831407	2234443	100.00%	19.309%
100	30.411	4648	4650	4651	rBV2	16542	12952	0.58%	0.112%

Sum of corrected areas: 11571972

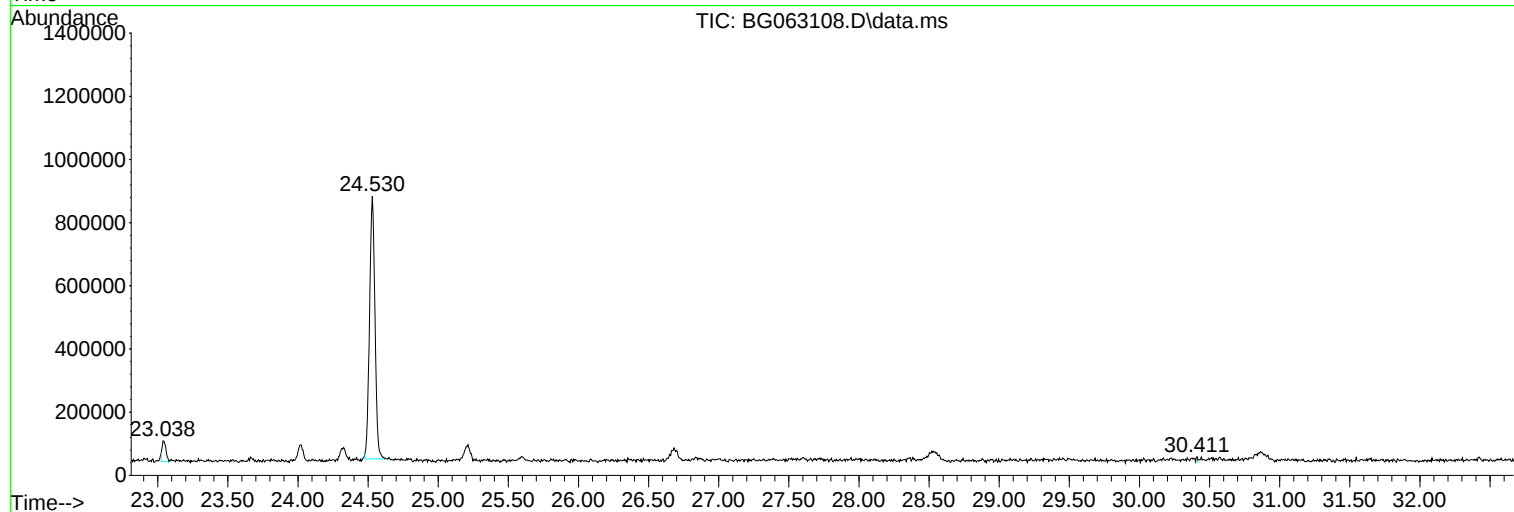
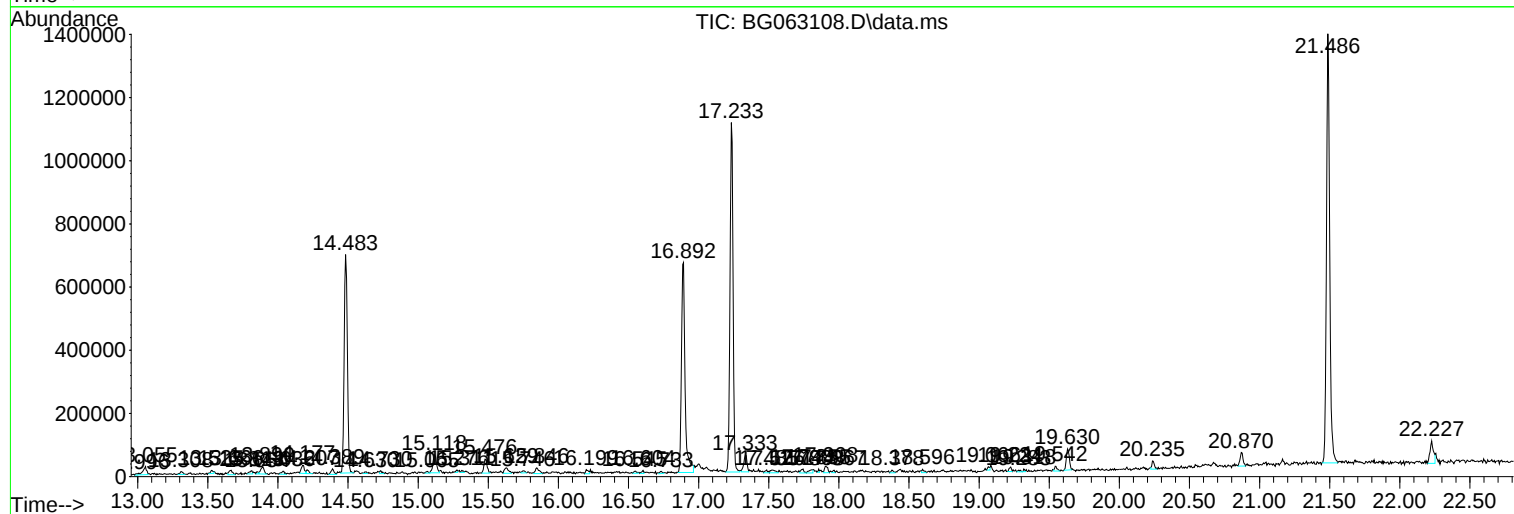
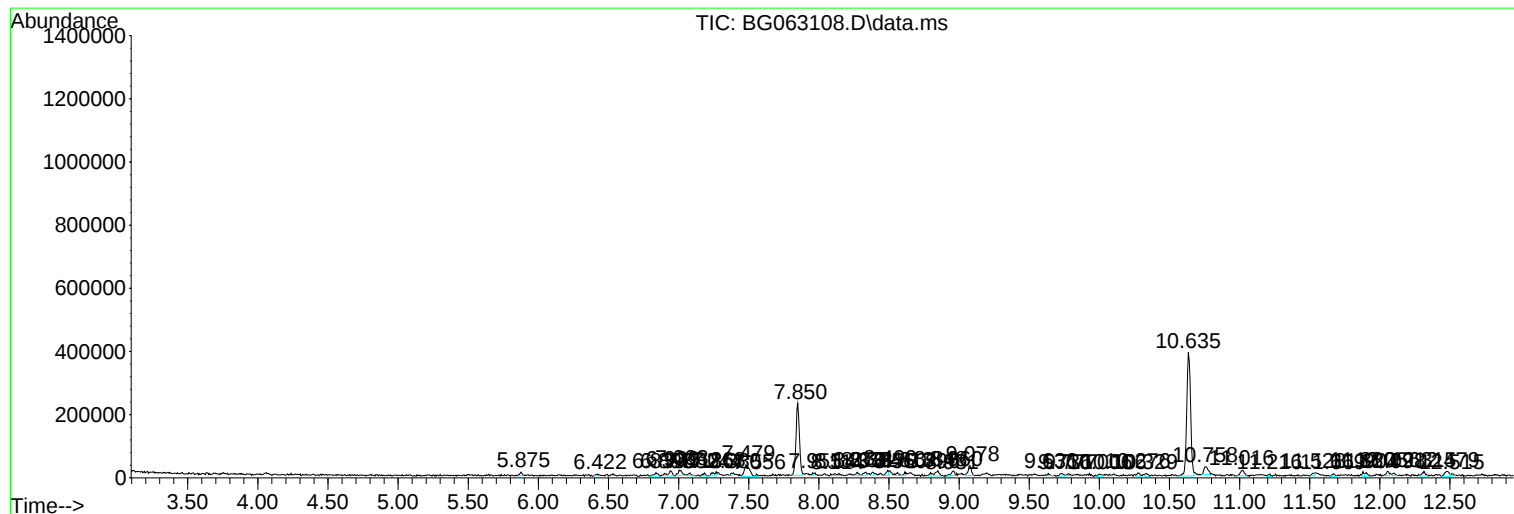
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P



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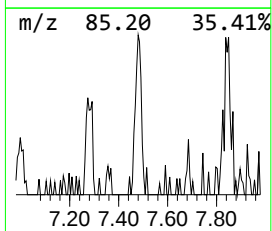
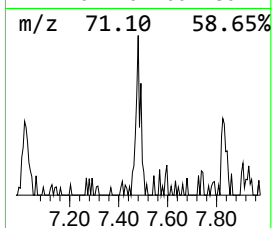
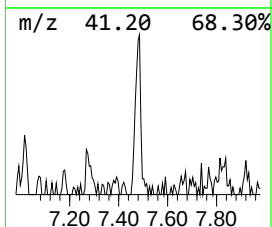
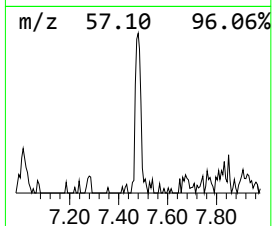
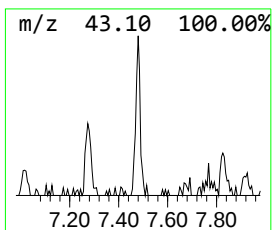
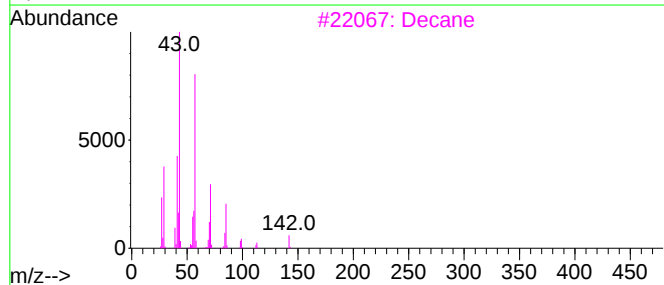
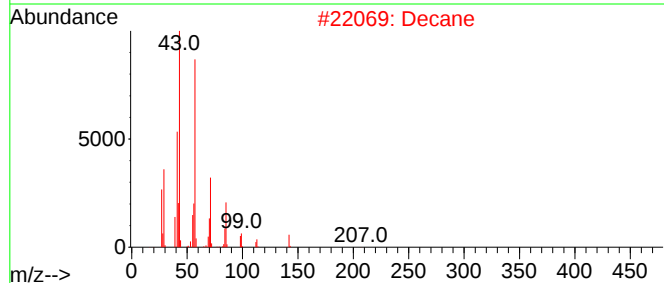
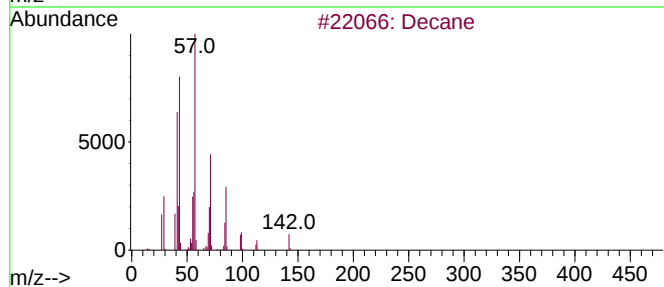
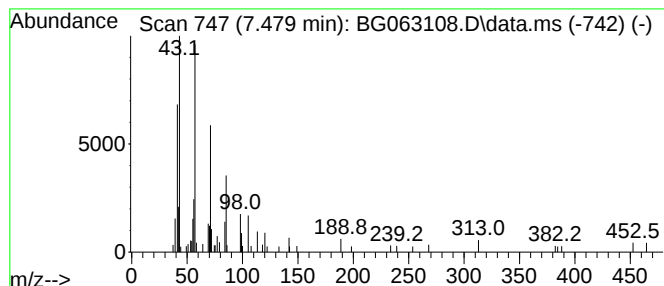
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.479	5.02 ng/ul	97669	1,4-Dichlorobenzene-d4			7.850
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	76
2	Decane		142	C10H22	000124-18-5	74
3	Decane		142	C10H22	000124-18-5	68
4	Nonadecane		268	C19H40	000629-92-5	58
5	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	50



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Instrument :
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ClientSampleId :
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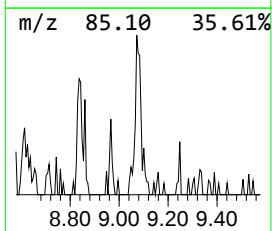
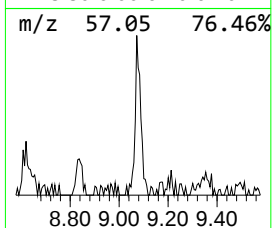
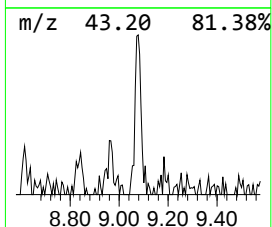
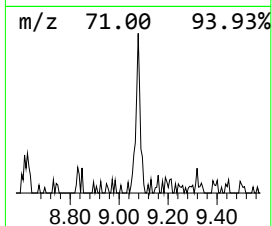
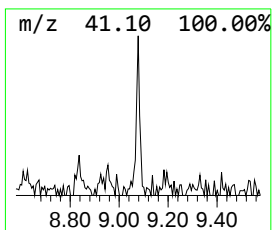
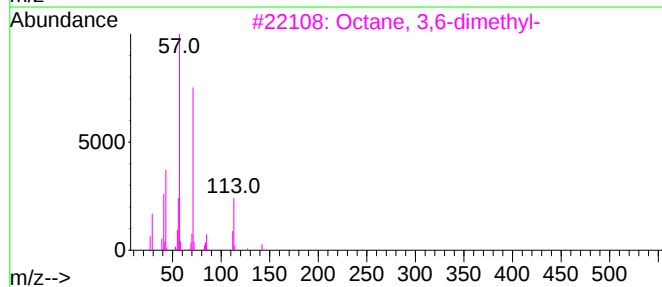
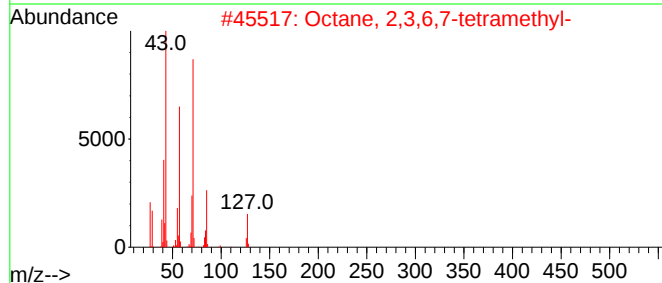
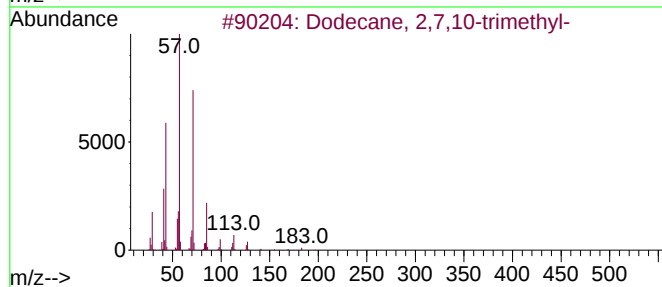
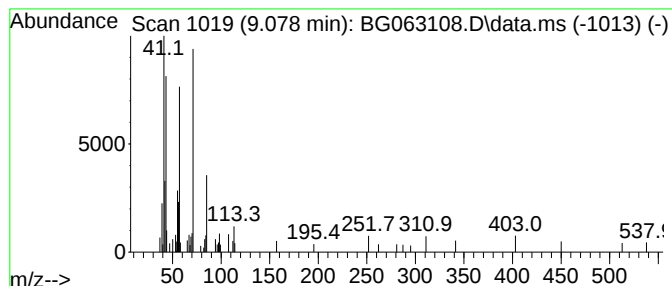
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.078	2.69 ng/ul	52386	1,4-Dichlorobenzene-d4	7.850

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	53
2			Octane, 2,3,6,7-tetramethyl-	170	C12H26	052670-34-5	53
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	47
4			3-Cyclopentylpropionamide, N-met...	195	C12H21NO	1000340-38-1	46
5			Dotriacontane	451	C32H66	000544-85-4	38



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA 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
(DEL) Alkane: S...	7.479	5.0	ng/u1	97669	1	7.850	389050	20.0
(DEL) Alkane: S...	9.078	2.7	ng/u1	52386	1	7.850	389050	20.0