LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022823\
Data File : BG056757.D
Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBZT7

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-BG022323.M
Title : SVOA CALIBRATION

Signal : TIC: BG056757.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.697	25	27	31	rBV2	2057	2044	0.52%	0.044%
2	4.149	101	104	109	rBV2	3758	5918	1.50%	0.128%
3	5.448	320	325	331	rVB3	8306	13318	3.37%	0.288%
4	7.539	673	681	687	rBV	40711	69732	17.64%	1.507%
5	7.721	706	712	718	rVB	28365	44903	11.36%	0.970%
6	7.874	734	738	743	rVV3	3278	5174	1.31%	0.112%
7	7.939	743	749	756	rVB2	35583	62124	15.71%	1.342%
8	8.421	824	831	839	rVB	56902	93477	23.65%	2.020%
9	8.732	877	884	888	rBV4	7097	10584	2.68%	0.229%
10	9.102	941	947	954	rBV2	43939	72840	18.43%	1.574%
11	9.243	966	971	981	rBV3	7043	11528	2.92%	0.249%
12	9.590	1024	1030	1037	rBV	39041	68375	17.30%	1.477%
13	9.972	1091	1095	1098	rVB3	1212	1664	0.42%	0.036%
14	10.318	1148	1154	1161	rVB2	34240	59483	15.05%	1.285%
15	10.859	1239	1246	1252	rBV2	52131	87382	22.10%	1.888%
16	11.018	1269	1273	1277	rBV2	1073	1768	0.45%	0.038%
17	11.070	1277	1282	1286	rVB2	13516	19313	4.89%	0.417%
18	11.253	1307	1313	1321	rVB	79606	143243	36.23%	3.095%
19	11.376	1327	1334	1341	rBV2	27325	51341	12.99%	1.109%
20	12.674	1550	1555	1558	rBV3	1810	1973	0.50%	0.043%
21	12.727	1560	1564	1570	rVB3	2641	3470	0.88%	0.075%
22	12.804	1575	1577	1583	rVB2	1394	2159	0.55%	0.047%
23	13.397	1674	1678	1684	rVB4	1520	2378	0.60%	0.051%
24	13.538	1696	1702	1709	rVV2	15032	21056	5.33%	0.455%
25	13.738	1731	1736	1741	rVB2	19429	25890	6.55%	0.559%
26	13.791	1741	1745	1749	rVV3	6746	10645	2.69%	0.230%
27	13.820	1749	1750	1753	rVB2	2787	2384	0.60%	0.052%
28	13.891	1757	1762	1766	rBV	1194	2059	0.52%	0.044%
29	14.226	1814	1819	1823	rBV3	11476	13879	3.51%	0.300%
30	14.349	1836	1840	1844	rVB2	4374	5162	1.31%	0.112%
31	14.408	1844	1850	1857	rBV	112672	160531	40.61%	3.469%
32	14.461	1857	1859	1862	rVV	2213	2619	0.66%	0.057%
33	14.508	1862	1867	1872	rVV	116264	162249	41.04%	3.506%
34	14.625	1883	1887	1889	rBV2	4791	6246	1.58%	0.135%
35	14.660	1889	1893	1897	rVV3	3988	4695	1.19%	0.101%
36	14.719	1897	1903	1910	rVB2	118133	180871	45.75%	3.908%

LSC Area Percent Report

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Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBZT7

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-BG022323.M
Title : SVOA CALIBRATION

37	14.878	1923	1930	1935	rBV2	2621	5010	1.27%	0.108%
38	14.989	1943	1949	1950	rBV	36658	46867	11.86%	1.013%
39	15.019	1950	1954	1961	rVB2	150854	233968	59.18%	5.056%
40	15.077	1961	1964	1969	rVB	1678	2463	0.62%	0.053%
41	15.171	1969	1980	1990	rBV	49019	78282	19.80%	1.692%
42	15.248	1990	1993	1998	rVB3	3384	3759	0.95%	0.081%
43	15.336	2003	2008	2010	rVB2	1712	2305	0.58%	0.050%
44	15.759	2072	2080	2087	rVV2	16009	22402	5.67%	0.484%
45	16.006	2115	2122	2129	rBV	160419	252385	63.84%	5.454%
46	16.100	2133	2138	2146	rBV	55564	79314	20.06%	1.714%
47	16.288	2165	2170	2173	rVB4	7670	10471	2.65%	0.226%
48	16.352	2176	2181	2186	rBV	52562	75459	19.09%	1.631%
49	16.582	2217	2220	2224	rVB4	4186	4478	1.13%	0.097%
50	16.629	2224	2228	2231	rBV3	1413	1886	0.48%	0.041%
51	16.711	2238	2242	2246	rBV4	1523	2260	0.57%	0.049%
52	17.204	2322	2326	2330	rVB5	1985	2520	0.64%	0.054%
53	17.246	2330	2333	2336	rBV2	2059	2519	0.64%	0.054%
54	17.422	2357	2363	2370	rVV4	12317	17649	4.46%	0.381%
55	17.686	2407	2408	2410	rVV	2033	1831	0.46%	0.040%
56	17.757	2413	2420	2424	rVV	202483	316403	80.04%	6.837%
57	17.798	2424	2427	2431	rVV2	16343	22353	5.65%	0.483%
58	17.851	2431	2436	2443	rVV2	183416	269701	68.22%	5.828%
59	17.909	2444	2446	2449	rVV3	4312	5873	1.49%	0.127%
60	17.939	2449	2451	2457	rVB3	5435	5436	1.38%	0.117%
61	17.992	2457	2460	2463	rBV3	1802	2070	0.52%	0.045%
62	18.144	2483	2486	2488	rBV	2921	3061	0.77%	0.066%
63	18.350	2517	2521	2524	rBV2	1025	1882	0.48%	0.041%
64	18.391	2524	2528	2532	rVB3	8329	9692	2.45%	0.209%
65	18.497	2542	2546	2550	rVB3	5625	7952	2.01%	0.172%
66	18.597	2559	2563	2571	rBV6	23207	46771	11.83%	1.011%
67	18.667	2571	2575	2582	rBV6	3843	6299	1.59%	0.136%
68	18.808	2596	2599	2600	rBV2	3975	3807	0.96%	0.082%
69	18.850	2603	2606	2611	rVB5	9487	11973	3.03%	0.259%
70	19.008	2629	2633	2639	rVB4	5169	7643	1.93%	0.165%
71	19.102	2645	2649	2650	rBV2	2278	2660	0.67%	0.057%
72	19.196	2663	2665	2668	rVB2	3252	2561	0.65%	0.055%
73	19.396	2695	2699	2703	rVV3	2518	3209	0.81%	0.069%
74	19.502	2714	2717	2721	rBV5	2437	3675	0.93%	0.079%
75	19.596	2731	2733	2735	rBV	1481	1957	0.50%	0.042%
76	19.654	2740	2743	2745	rBV3	2377	3070	0.78%	0.066%

LSC Area Percent Report

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Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

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
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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Title : SVOA CALIBRATION

77	19.778	2760	2764	2769	rVB	31121	44817	11.34%	0.968%
78	19.848	2772	2776	2777	rBV2	4690	5412	1.37%	0.117%
79	19.931	2787	2790	2792	rBV3	3412	4698	1.19%	0.102%
80	20.019	2792	2805	2806	rBV3	16586	48785	12.34%	1.054%
81	20.113	2815	2821	2831	rVB	242302	395323	100.00%	8.542%
82	20.318	2852	2856	2861	rVB6	3485	5252	1.33%	0.113%
83	20.454	2874	2879	2880	rBV5	3586	5100	1.29%	0.110%
84	20.501	2885	2887	2890	rBV3	2048	3082	0.78%	0.067%
85	20.536	2890	2893	2897	rVV5	5187	6556	1.66%	0.142%
86	20.595	2900	2903	2906	rVB3	4889	7044	1.78%	0.152%
87	20.665	2913	2915	2918	rBV2	2012	2417	0.61%	0.052%
88	20.853	2943	2947	2949	rBV	10683	12547	3.17%	0.271%
89	21.100	2987	2989	2992	rVB3	6821	7269	1.84%	0.157%
90	21.370	3033	3035	3038	rBV2	3206	3117	0.79%	0.067%
91	21.588	3067	3072	3075	rBV7	6722	10117	2.56%	0.219%
92	21.640	3075	3081	3089	rVV3	38693	70647	17.87%	1.527%
93	22.063	3145	3153	3159	rBV	191858	350529	88.67%	7.574%
94	22.839	3283	3285	3287	rBV3	3312	2848	0.72%	0.062%
95	23.861	3455	3459	3462	rBV5	7364	10369	2.62%	0.224%
96	25.254	3694	3696	3698	rBV2	3780	2478	0.63%	0.054%
97	25.336	3702	3710	3719	rBV3	103374	289248	73.17%	6.250%
98	25.589	3743	3753	3763	rVB2	105852	313981	79.42%	6.785%
99	26.975	3983	3989	3991	rBV7	3407	7639	1.93%	0.165%
100	31.071	4684	4686	4689	rBV3	2608	2139	0.54%	0.046%

Sum of corrected areas: 4627767

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Instrument :
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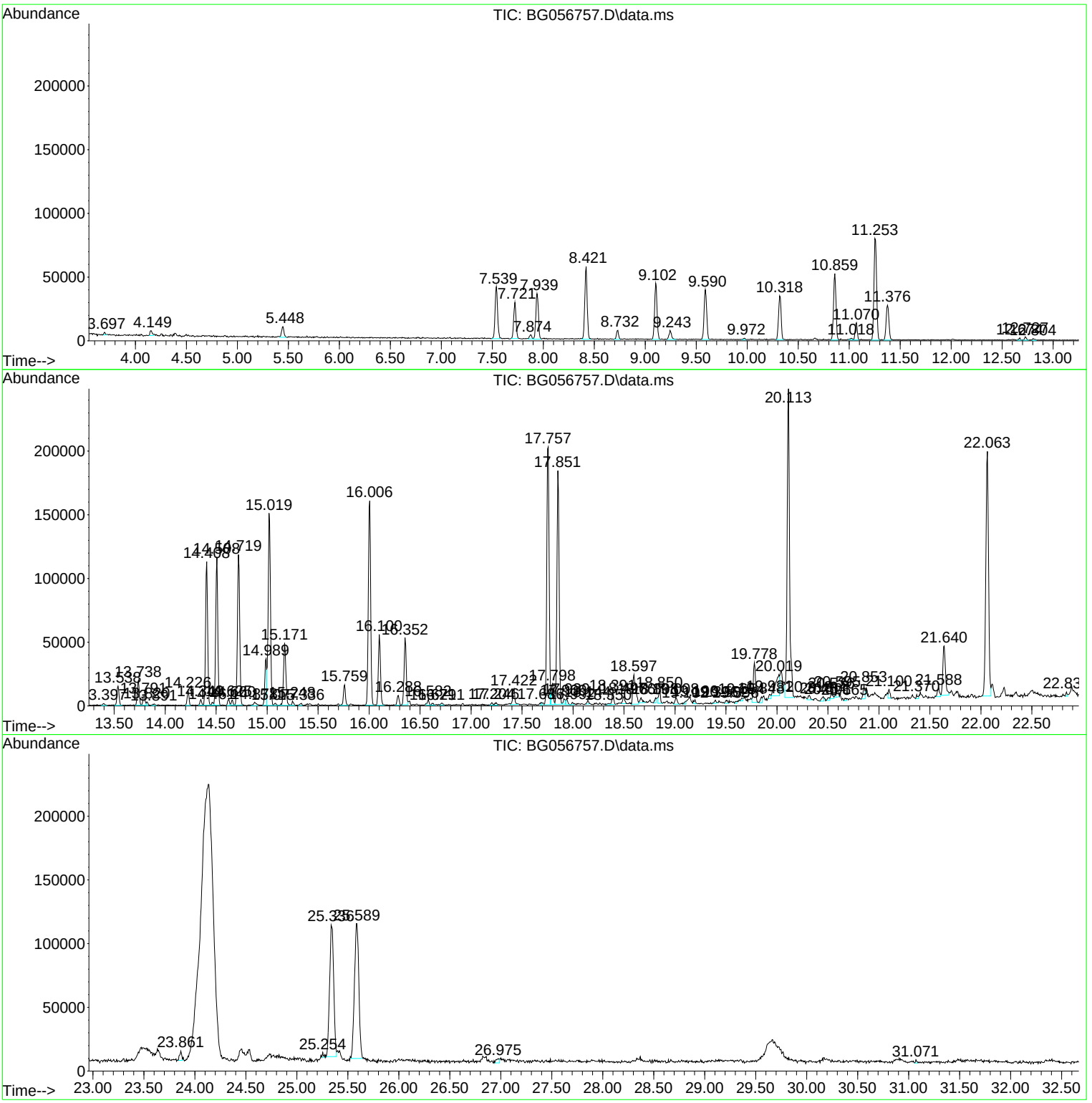
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022823\
Data File : BG056757.D
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Sample : 01648-09
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ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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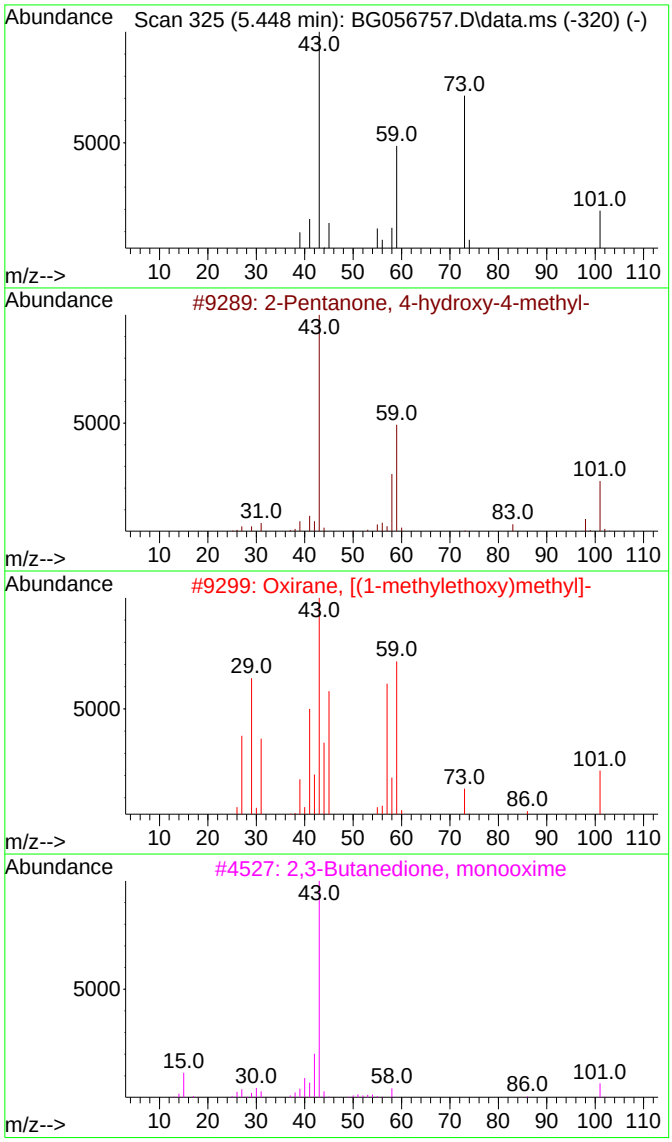
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

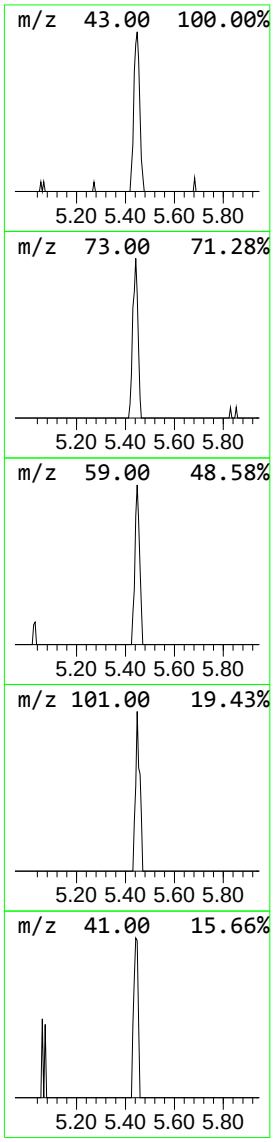
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.448	2.85 ng/ul	13318	1,4-Dichlorobenzene-d4	8.421

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	9
2	Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	9
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
4	Diacetamide	101	C4H7NO2	000625-77-4	7
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022823\
Data File : BG056757.D
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Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

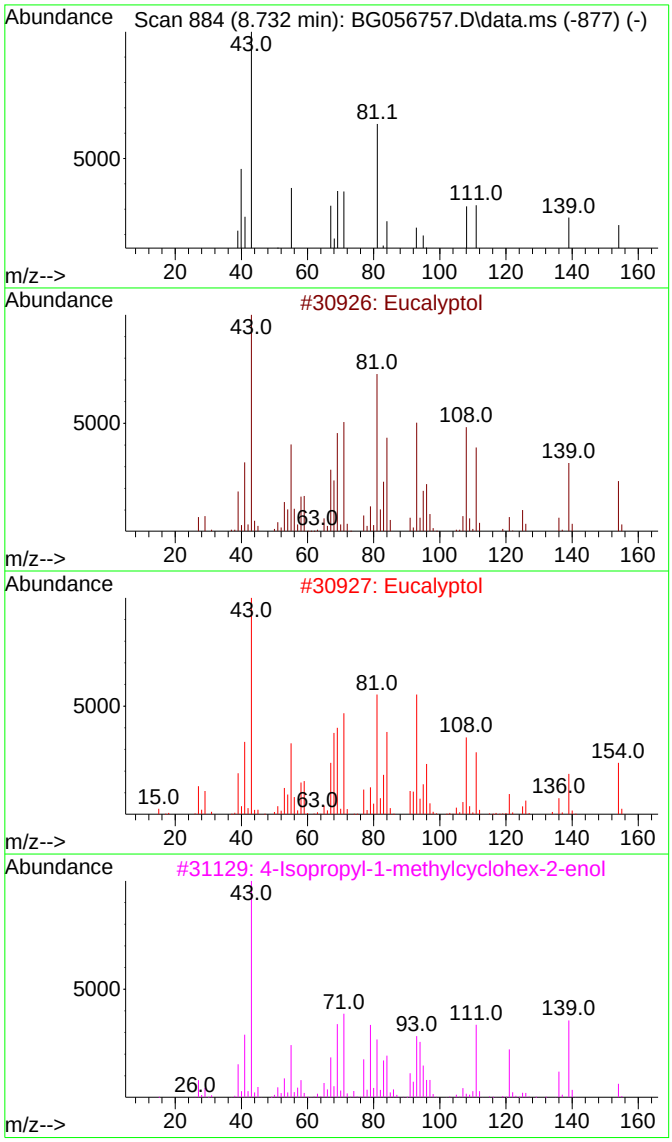
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

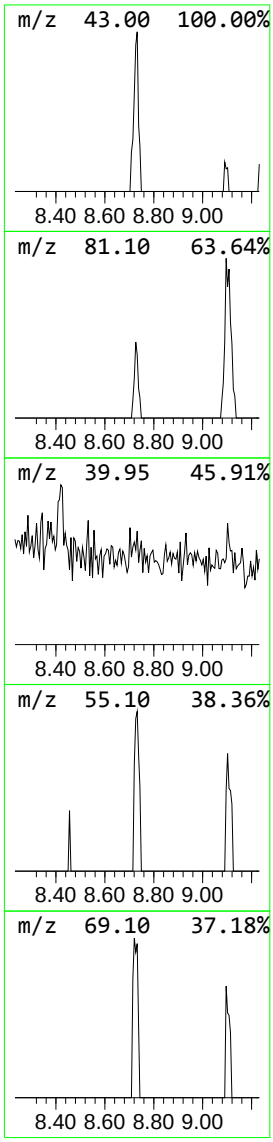
****

Peak Number 2 Eucalyptol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.732	2.26 ng/ul	10584	1,4-Dichlorobenzene-d4	8.421

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	87
2	Eucalyptol	154	C10H18O	000470-82-6	40
3	4-Isopropyl-1-methylcyclohex-2-enol	154	C10H18O	000619-62-5	37
4	2-Cyclohexen-1-ol, 1-methyl-4-(1...	154	C10H18O	029803-82-5	35
5	1-Cyclopentene-1-methanol, .alph...	154	C10H18O	056666-68-3	25





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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

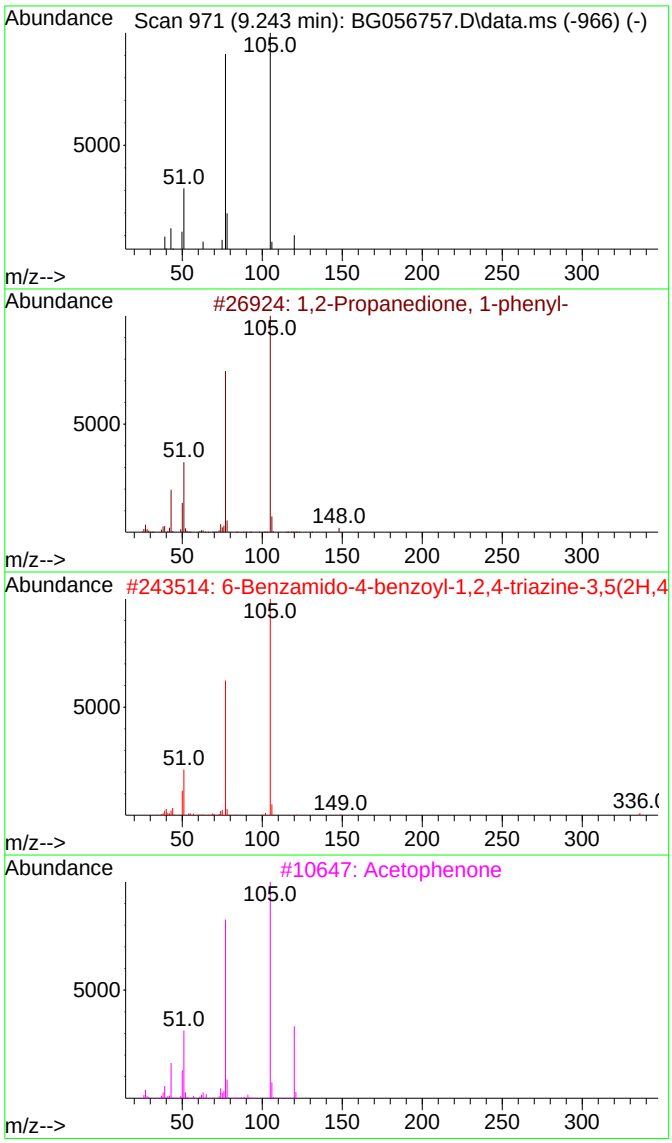
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

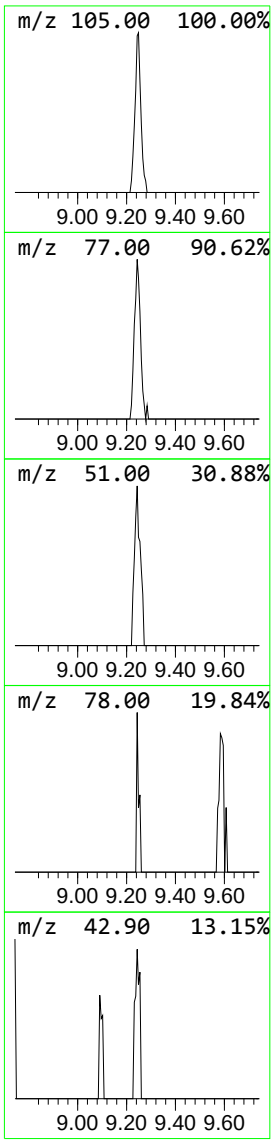
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1,2-Propanedione, 1-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.243	2.47 ng/ul	11528	1,4-Dichlorobenzene-d4	8.421

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	50
2	6-Benzamido-4-benzoyl-1,2,4-tria...	336	C17H12N4O4	1000224-14-0	50
3	Acetophenone	120	C8H8O	000098-86-2	40
4	Acetophenone	120	C8H8O	000098-86-2	40
5	Acetophenone	120	C8H8O	000098-86-2	40





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Instrument :
BNA_G
ClientSampleId :
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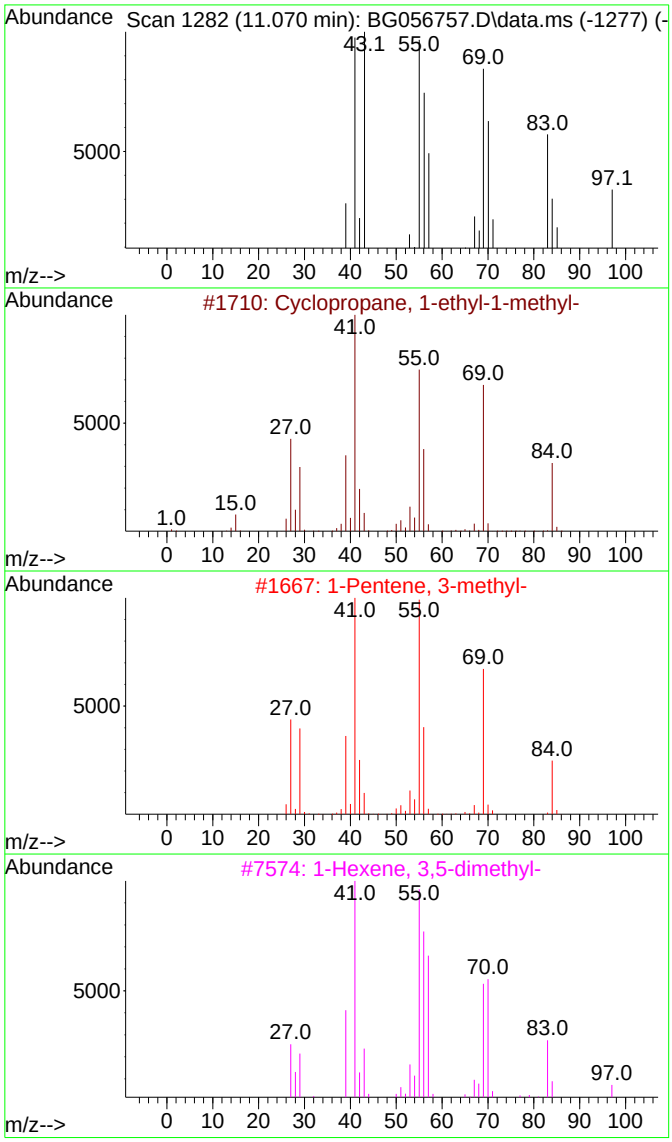
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

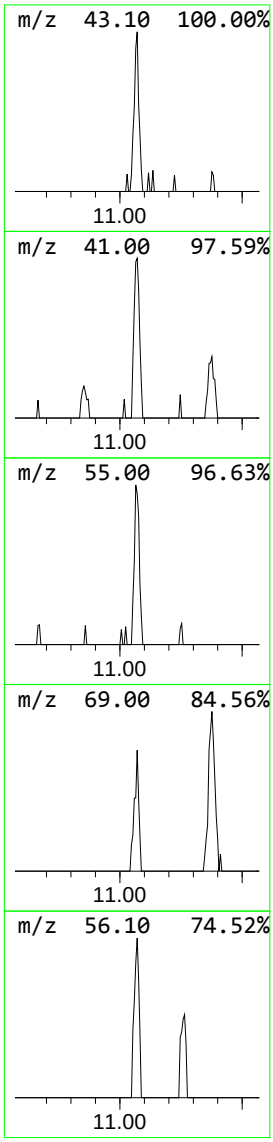
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Cyclic11.070 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.070	2.70 ng/ul	19313	Naphthalene-d8	11.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	43
2	1-Pentene, 3-methyl-	84	C6H12	000760-20-3	38
3	1-Hexene, 3,5-dimethyl-	112	C8H16	007423-69-0	38
4	1-Octanol	130	C8H18O	000111-87-5	37
5	Pentane, 2-cyclopropyl-	112	C8H16	005458-16-2	35





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Data File : BG056757.D
Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

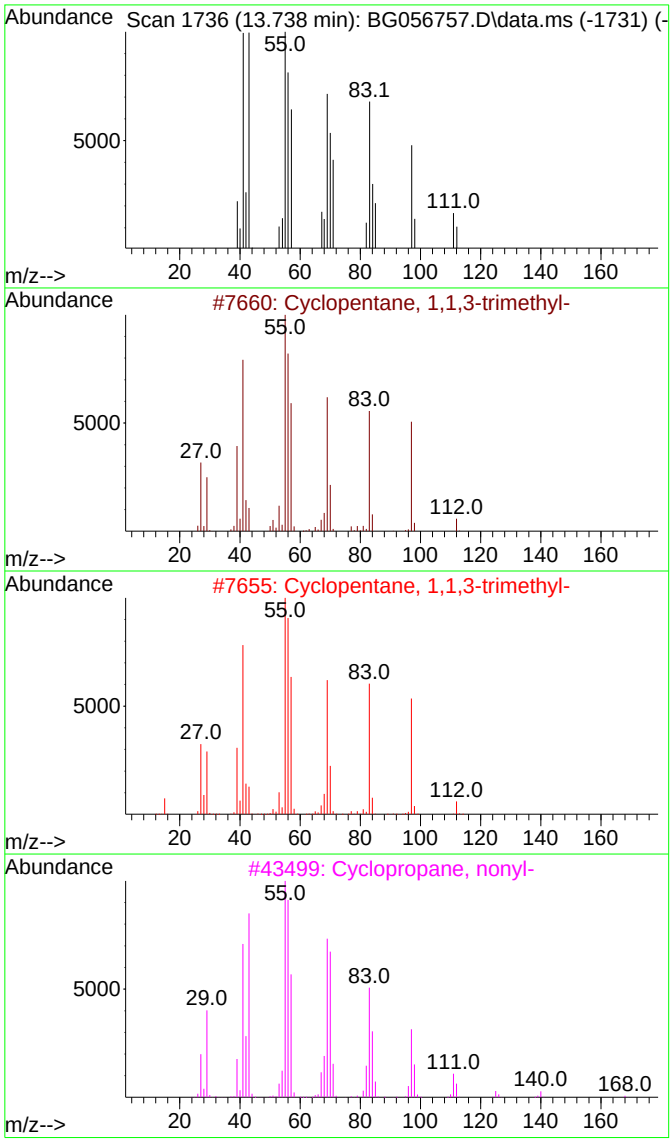
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

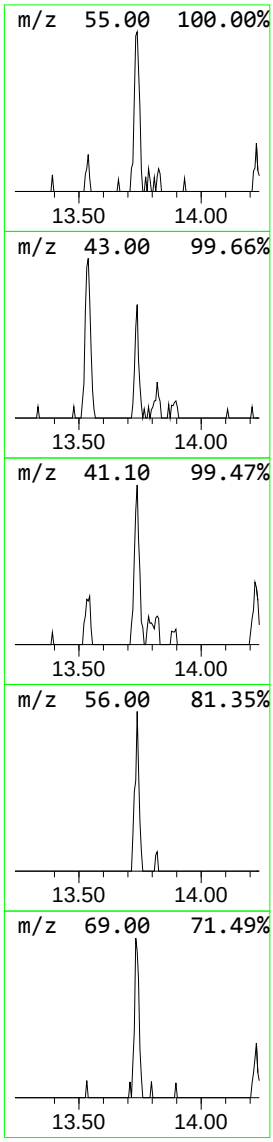
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic13.738 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.738	2.21 ng/ul	25890	Acenaphthene-d10	15.019

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	64
2	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	64
3	Cyclopropane, nonyl-	168	C12H24	074663-85-7	52
4	Cyclopropane, octyl-	154	C11H22	001472-09-9	50
5	1-Decanol	158	C10H22O	000112-30-1	47





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Data File : BG056757.D
Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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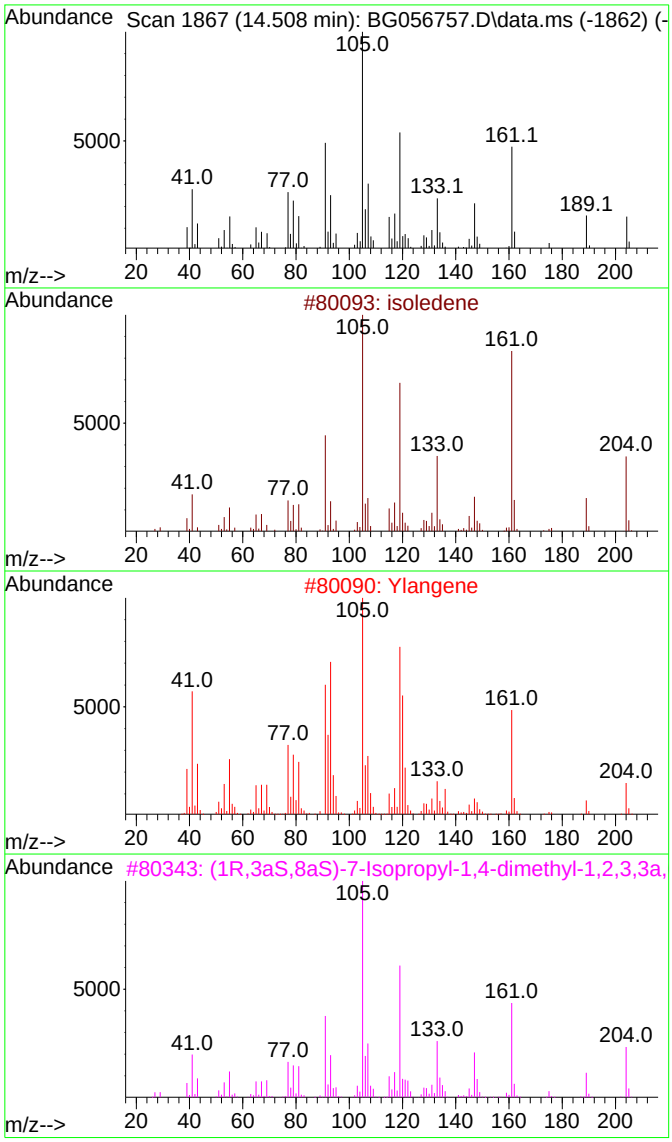
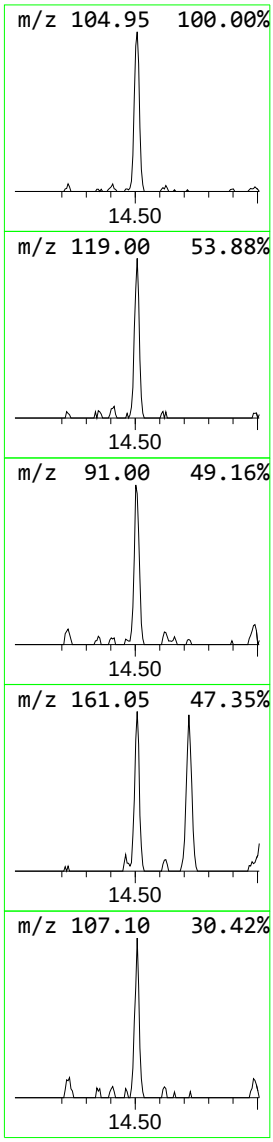
Peak Number 6 isoledene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.508	13.87 ng/ul	162249	Acenaphthene-d10	15.019

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	isoledene	204	C15H24	095910-36-4	91
2	Ylangene	204	C15H24	014912-44-8	64
3	(1R,3aS,8aS)-7-Isopropyl-1,4-dim...	204	C15H24	036577-33-0	58
4	(-)-Aristolene	204	C15H24	006831-16-9	55
5	(-)-Aristolene	204	C15H24	006831-16-9	50





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Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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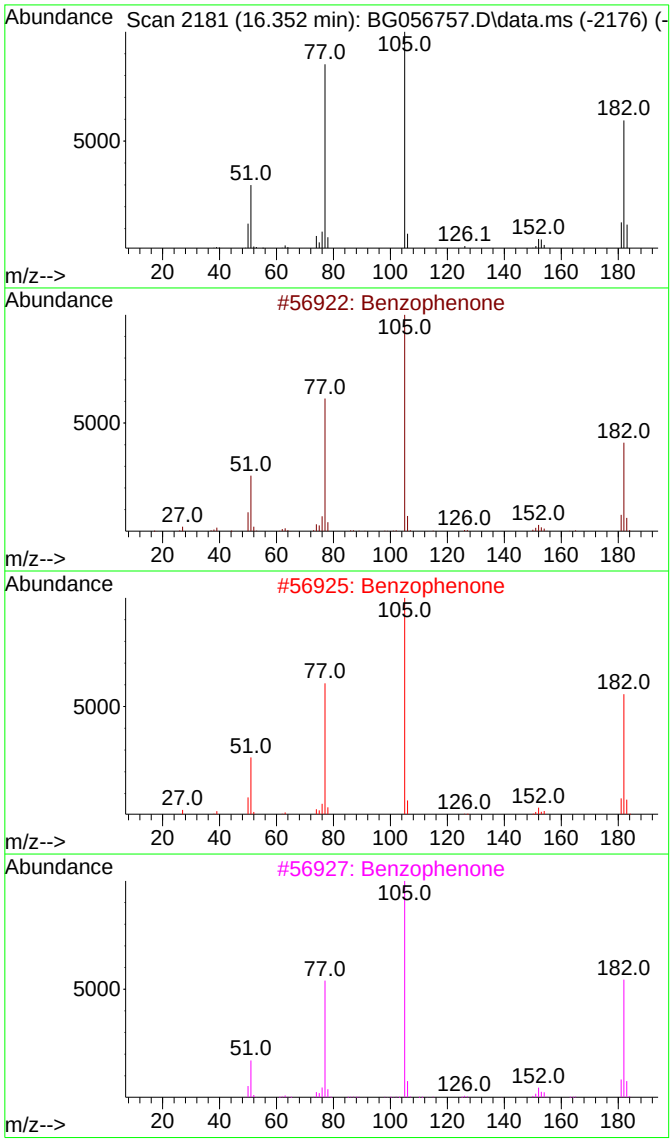
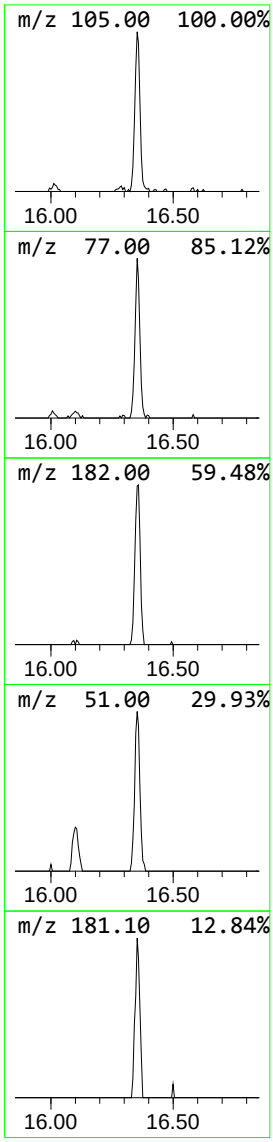
Peak Number 7 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.352	6.45 ng/ul	75459	Acenaphthene-d10	15.019

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	89
5	Benzophenone	182	C13H10O	000119-61-9	80





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Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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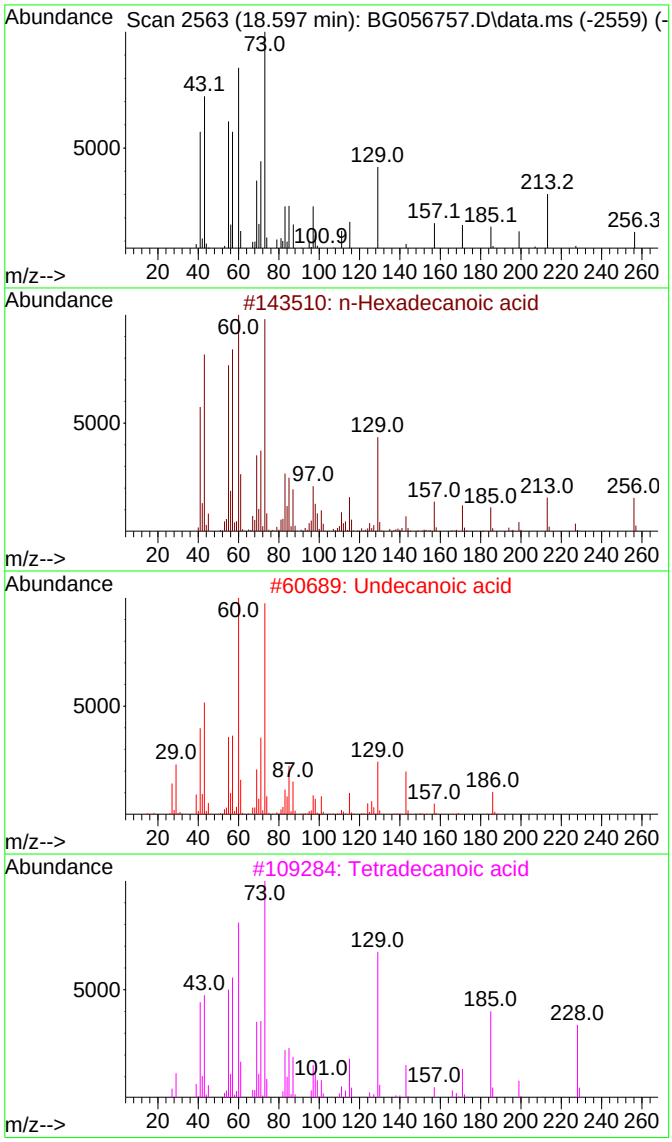
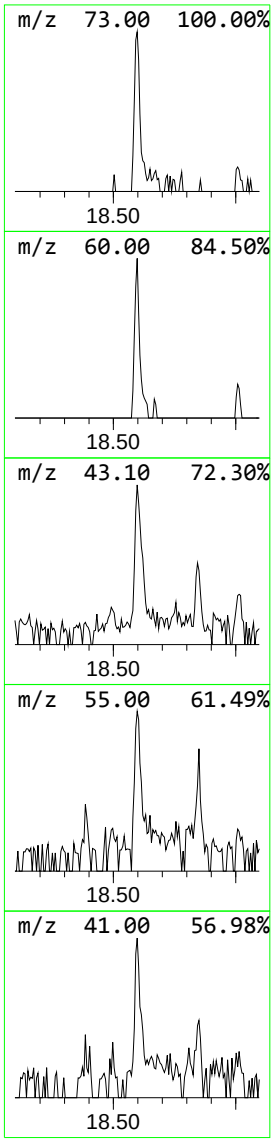
Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.597	2.96 ng/ul	46771	Phenanthrene-d10	17.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2	Undecanoic acid	186	C11H22O2	000112-37-8	81
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4	n-Decanoic acid	172	C10H20O2	000334-48-5	53
5	Tridecanoic acid	214	C13H26O2	000638-53-9	49





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Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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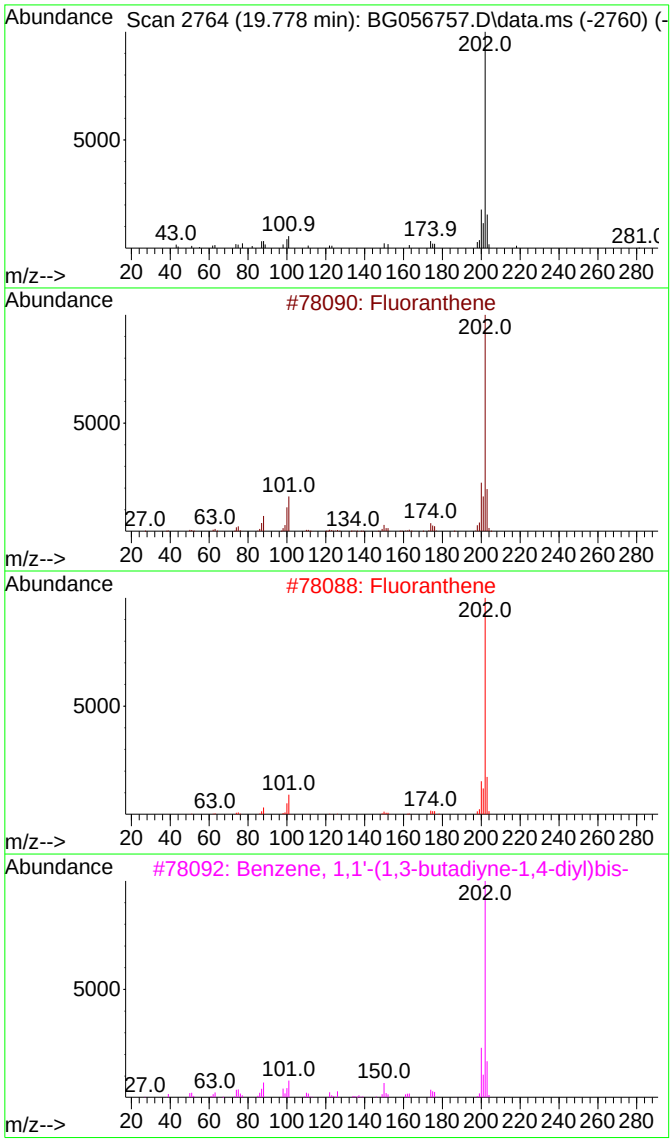
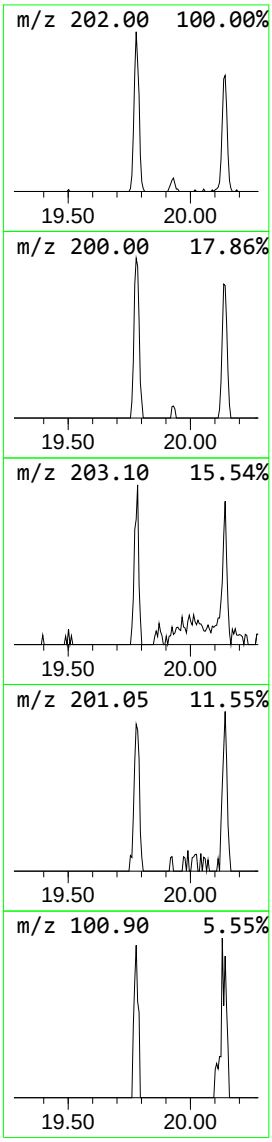
Peak Number 9 Fluoran thene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.778	2.83 ng/ul	44817	Phenanthrene-d10	17.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene	202	C16H10	000206-44-0	90
2	Fluoranthene	202	C16H10	000206-44-0	86
3	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	83
4	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	83
5	Fluoranthene	202	C16H10	000206-44-0	81





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Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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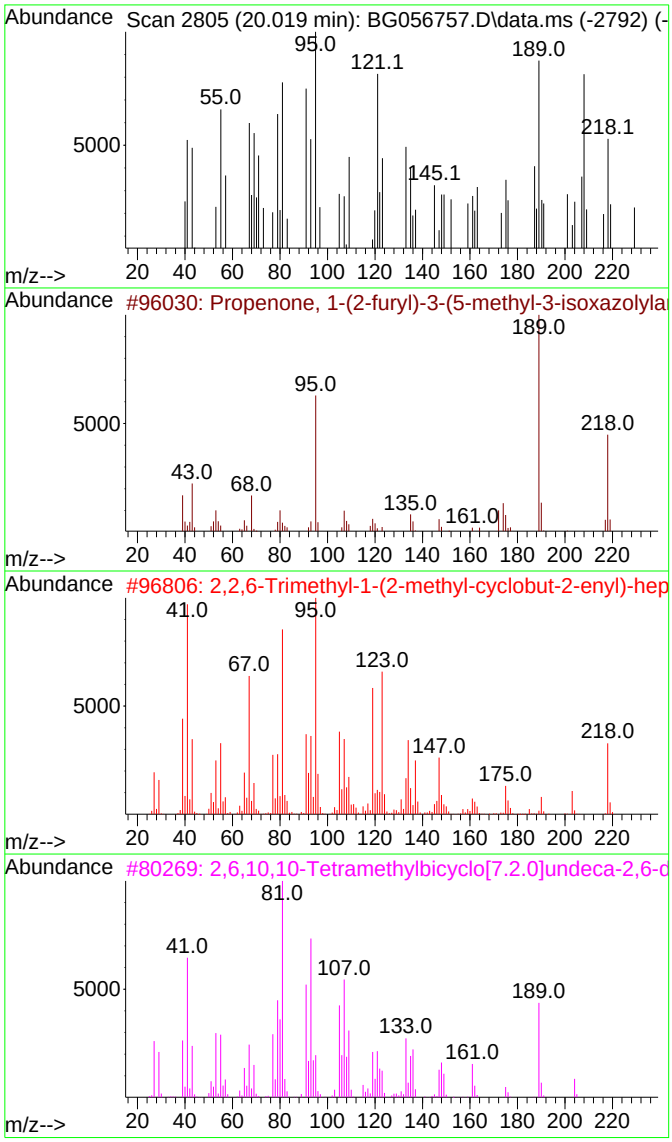
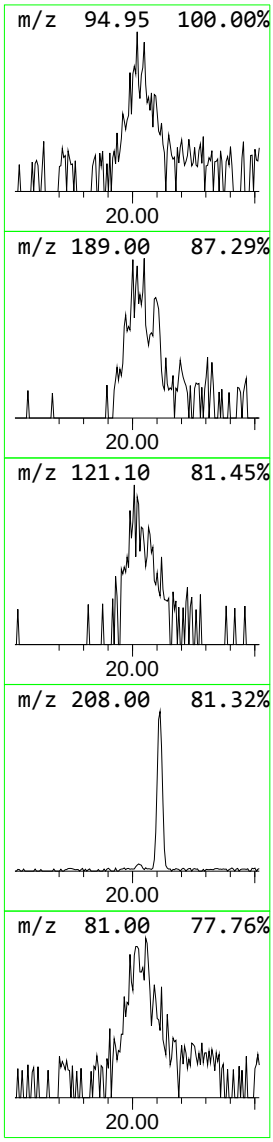
Peak Number 10 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.019	2.78 ng/ul	48785	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propenone, 1-(2-furyl)-3-(5-meth...	218	C11H10N2O3	186957-33-5	27
2	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	25
3	2,6,10,10-Tetramethylbicyclo[7.2...	204	C15H24	136296-37-2	22
4	Sesquirosefuran	218	C15H22O	039007-93-7	18
5	Pyrene, hexadecahydro-	218	C16H26	002435-85-0	18





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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

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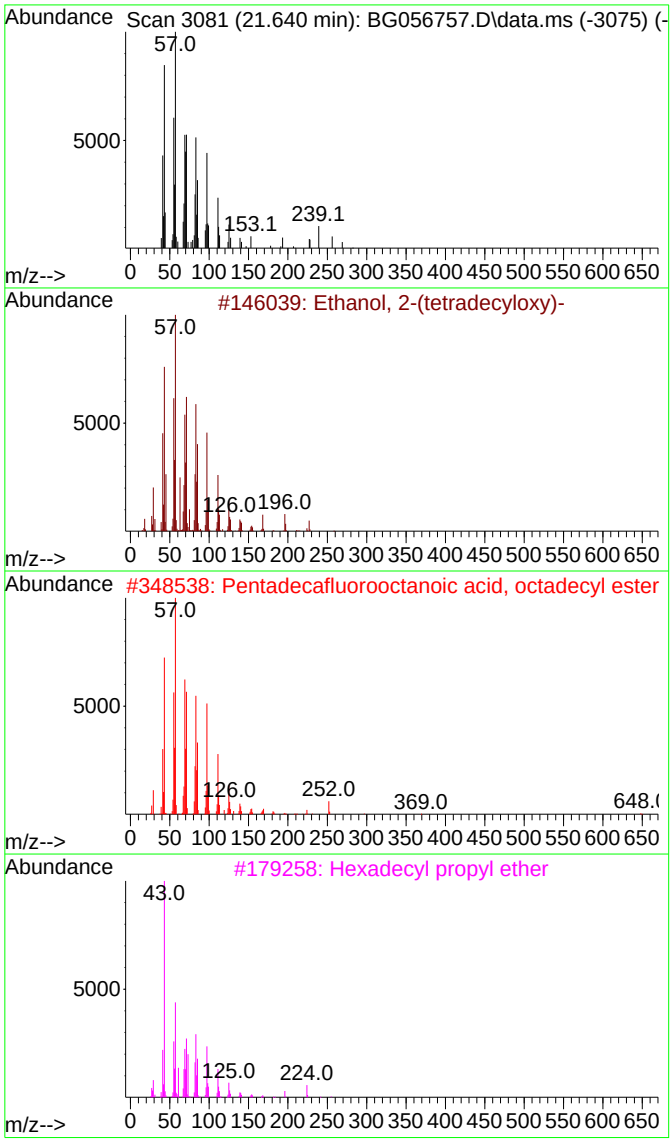
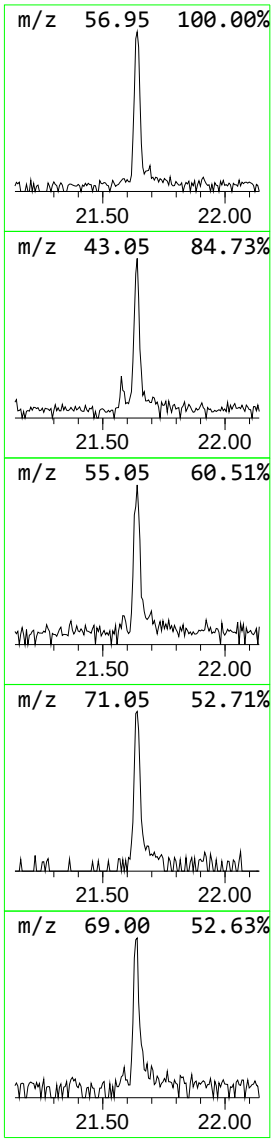
Peak Number 11 Ethanol, 2-(tetradecyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.640	4.03 ng/ul	70647	Chrysene-d12	22.063

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	87
3	Hexadecyl propyl ether	284	C19H40O	1000406-27-9	83
4	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	81
5	Propyl triacontyl ether	481	C33H68O	1000406-28-6	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022823\
Data File : BG056757.D
Acq On : 28 Feb 2023 16:21
Operator : CG/JU
Sample : 01648-09
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBZT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	5.448	2.9	ng/ul	13318	1	8.421	93477	20.0
Eucalyptol	8.732	2.3	ng/ul	10584	1	8.421	93477	20.0
1,2-Propanedion...	9.243	2.5	ng/ul	11528	1	8.421	93477	20.0
(DEL) Alkane: C...	11.070	2.7	ng/ul	19313	2	11.253	143243	20.0
(DEL) Alkane: C...	13.738	2.2	ng/ul	25890	3	15.019	233968	20.0
isoledene	14.508	13.9	ng/ul	162249	3	15.019	233968	20.0
Benzophenone	16.352	6.5	ng/ul	75459	3	15.019	233968	20.0
n-Hexadecanoic ...	18.597	3.0	ng/ul	46771	4	17.757	316403	20.0
Fluoran thene	19.778	2.8	ng/ul	44817	4	17.757	316403	20.0
unknown-01	20.019	2.8	ng/ul	48785	5	22.063	350529	20.0
Ethanol, 2-(tet...	21.640	4.0	ng/ul	70647	5	22.063	350529	20.0