

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
 Data File : BG047559.D
 Acq On : 4 Dec 2020 5:32
 Operator : CG/JU
 Sample : L4855-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CB7Y9

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.490	23	26	29	rBV4	3519	4961	0.40%	0.030%
2	3.596	39	44	51	rVB4	9104	14908	1.20%	0.090%
3	4.007	113	114	117	rBV	2970	3377	0.27%	0.020%
4	4.371	169	176	184	rBV3	12157	27433	2.21%	0.165%
5	4.735	234	238	245	rVB4	14274	24886	2.00%	0.150%
6	5.006	280	284	288	rBV	2246	3214	0.26%	0.019%
7	5.082	292	297	300	rVB	2485	2936	0.24%	0.018%
8	5.523	367	372	376	rBV	2671	3933	0.32%	0.024%
9	5.864	424	430	433	rVV	1956	4557	0.37%	0.027%
10	6.204	484	488	489	rBV	3920	4352	0.35%	0.026%
11	6.269	496	499	504	rVB3	7691	10681	0.86%	0.064%
12	7.109	638	642	644	rVB	2258	2781	0.22%	0.017%
13	7.168	648	652	653	rVV	2341	2812	0.23%	0.017%
14	7.191	654	656	660	rVB	3062	3022	0.24%	0.018%
15	7.332	675	680	681	rBV	3391	4864	0.39%	0.029%
16	7.391	683	690	700	rVB	227794	412380	33.20%	2.482%
17	7.561	712	719	729	rVB	129899	225448	18.15%	1.357%
18	7.779	748	756	765	rBV2	227154	402585	32.41%	2.423%
19	7.867	767	771	774	rVV4	3358	5386	0.43%	0.032%
20	7.896	774	776	778	rVV	2174	2699	0.22%	0.016%
21	7.926	778	781	786	rVB3	7482	9006	0.72%	0.054%
22	8.255	828	837	844	rBV	147661	268025	21.58%	1.613%
23	8.842	933	937	941	rVV2	4956	5482	0.44%	0.033%
24	8.948	945	955	964	rVV	265065	484244	38.98%	2.915%
25	9.060	971	974	977	rBV2	2054	2789	0.22%	0.017%
26	9.359	1020	1025	1028	rBV2	5052	7626	0.61%	0.046%
27	9.418	1028	1035	1042	rVV	216289	393783	31.70%	2.370%
28	9.477	1044	1045	1050	rVB	4184	4645	0.37%	0.028%
29	9.571	1056	1061	1066	rBV2	26353	39869	3.21%	0.240%
30	9.829	1098	1105	1109	rBV3	11326	20193	1.63%	0.122%
31	10.147	1152	1159	1166	rVB	221369	390476	31.43%	2.351%
32	10.464	1206	1213	1217	rBV4	4929	7560	0.61%	0.046%
33	10.687	1243	1251	1262	rBV	302578	547908	44.11%	3.298%
34	11.081	1311	1318	1326	rVV	184533	342598	27.58%	2.062%

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35	11.140	1326	1328	1332	rVV3	6800	8265	0.67%	0.050%
36	11.198	1332	1338	1348	rVB	236788	429737	34.59%	2.587%
37	11.809	1439	1442	1444	rBV	1917	2757	0.22%	0.017%
38	11.927	1459	1462	1465	rBV	1828	2945	0.24%	0.018%
39	11.992	1469	1473	1479	rVB2	5182	9853	0.79%	0.059%
40	12.074	1486	1487	1489	rBV	3150	2876	0.23%	0.017%
41	12.462	1550	1553	1557	rVB2	5475	7464	0.60%	0.045%
42	12.585	1570	1574	1576	rBV2	4636	6590	0.53%	0.040%
43	12.644	1579	1584	1588	rVV	4726	9946	0.80%	0.060%
44	12.708	1590	1595	1602	rVV2	6548	10305	0.83%	0.062%
45	12.761	1602	1604	1608	rVB	2594	3428	0.28%	0.021%
46	12.938	1628	1634	1637	rVB2	3453	6036	0.49%	0.036%
47	13.308	1694	1697	1701	rBV3	4784	7176	0.58%	0.043%
48	13.490	1724	1728	1734	rVB3	3033	6978	0.56%	0.042%
49	13.549	1734	1738	1739	rBV3	2744	2815	0.23%	0.017%
50	13.684	1756	1761	1763	rBV2	4732	6673	0.54%	0.040%
51	14.189	1844	1847	1851	rBV2	2627	4045	0.33%	0.024%
52	14.260	1852	1859	1865	rBV	502620	733745	59.06%	4.417%
53	14.436	1887	1889	1891	rVB	4005	3107	0.25%	0.019%
54	14.565	1903	1911	1918	rVB	497255	820343	66.04%	4.938%
55	14.706	1932	1935	1937	rBV2	4290	3886	0.31%	0.023%
56	14.741	1937	1941	1943	rVB3	5382	6996	0.56%	0.042%
57	14.865	1953	1962	1970	rBV2	301403	491736	39.58%	2.960%
58	15.029	1984	1990	1997	rBV	280174	457381	36.82%	2.753%
59	15.423	2054	2057	2059	rVB2	2696	2647	0.21%	0.016%
60	15.640	2091	2094	2096	rVV4	5727	7334	0.59%	0.044%
61	15.681	2099	2101	2105	rVB2	4935	6199	0.50%	0.037%
62	15.852	2123	2130	2137	rVV	683287	1016821	81.85%	6.121%
63	15.952	2140	2147	2155	rBV2	232239	354443	28.53%	2.134%
64	16.093	2165	2171	2176	rVB5	10522	17892	1.44%	0.108%
65	16.187	2183	2187	2188	rBV2	4206	3161	0.25%	0.019%
66	16.398	2219	2223	2225	rBV3	8484	11236	0.90%	0.068%
67	16.539	2245	2247	2250	rVV2	4301	5130	0.41%	0.031%
68	16.621	2259	2261	2264	rBV2	4627	5512	0.44%	0.033%
69	16.663	2266	2268	2272	rVB3	3122	3564	0.29%	0.021%
70	16.721	2276	2278	2279	rVV2	4686	3081	0.25%	0.019%
71	16.827	2292	2296	2298	rBV3	4965	4928	0.40%	0.030%

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72	16.862	2301	2302	2304	rBV2	3626	3743	0.30%	0.023%
73	16.898	2306	2308	2313	rVB6	5227	6660	0.54%	0.040%
74	17.050	2331	2334	2336	rBV3	3636	4060	0.33%	0.024%
75	17.139	2345	2349	2351	rBV4	5280	5858	0.47%	0.035%
76	17.174	2351	2355	2357	rBV5	4281	5359	0.43%	0.032%
77	17.274	2371	2372	2375	rVB3	4333	3432	0.28%	0.021%
78	17.332	2377	2382	2387	rBV6	8759	15270	1.23%	0.092%
79	17.397	2387	2393	2395	rBV4	4894	8857	0.71%	0.053%
80	17.415	2395	2396	2398	rVV2	6031	3299	0.27%	0.020%
81	17.491	2408	2409	2414	rVB5	7023	7727	0.62%	0.047%
82	17.532	2414	2416	2418	rBV3	3396	3809	0.31%	0.023%
83	17.603	2422	2428	2433	rBV	396690	607486	48.90%	3.657%
84	17.703	2439	2445	2454	rVB	663991	1071267	86.23%	6.449%
85	17.826	2464	2466	2468	rVB2	5326	3166	0.25%	0.019%
86	17.861	2468	2472	2473	rBV4	7962	9732	0.78%	0.059%
87	18.126	2515	2517	2518	rBV2	4959	4065	0.33%	0.024%
88	18.208	2529	2531	2532	rBV2	5346	4073	0.33%	0.025%
89	19.465	2742	2745	2748	rVB4	15961	21104	1.70%	0.127%
90	19.630	2771	2773	2780	rVB2	48178	67781	5.46%	0.408%
91	19.965	2825	2830	2840	rVB2	801075	1242273	100.00%	7.478%
92	21.486	3085	3089	3095	rVB2	156143	236112	19.01%	1.421%
93	21.880	3151	3156	3162	rVB	334200	574660	46.26%	3.459%
94	22.720	3293	3299	3306	rVB4	121989	280351	22.57%	1.688%
95	23.801	3477	3483	3491	rVB2	252832	513705	41.35%	3.092%
96	24.295	3561	3567	3572	rBV2	258805	589991	47.49%	3.552%
97	24.354	3572	3577	3588	rVB3	333584	770236	62.00%	4.637%
98	25.041	3685	3694	3706	rVB	373573	1007640	81.11%	6.066%
99	25.276	3727	3734	3745	rVB2	211220	648495	52.20%	3.904%
100	26.510	3939	3944	3956	rVB3	255345	713168	57.41%	4.293%

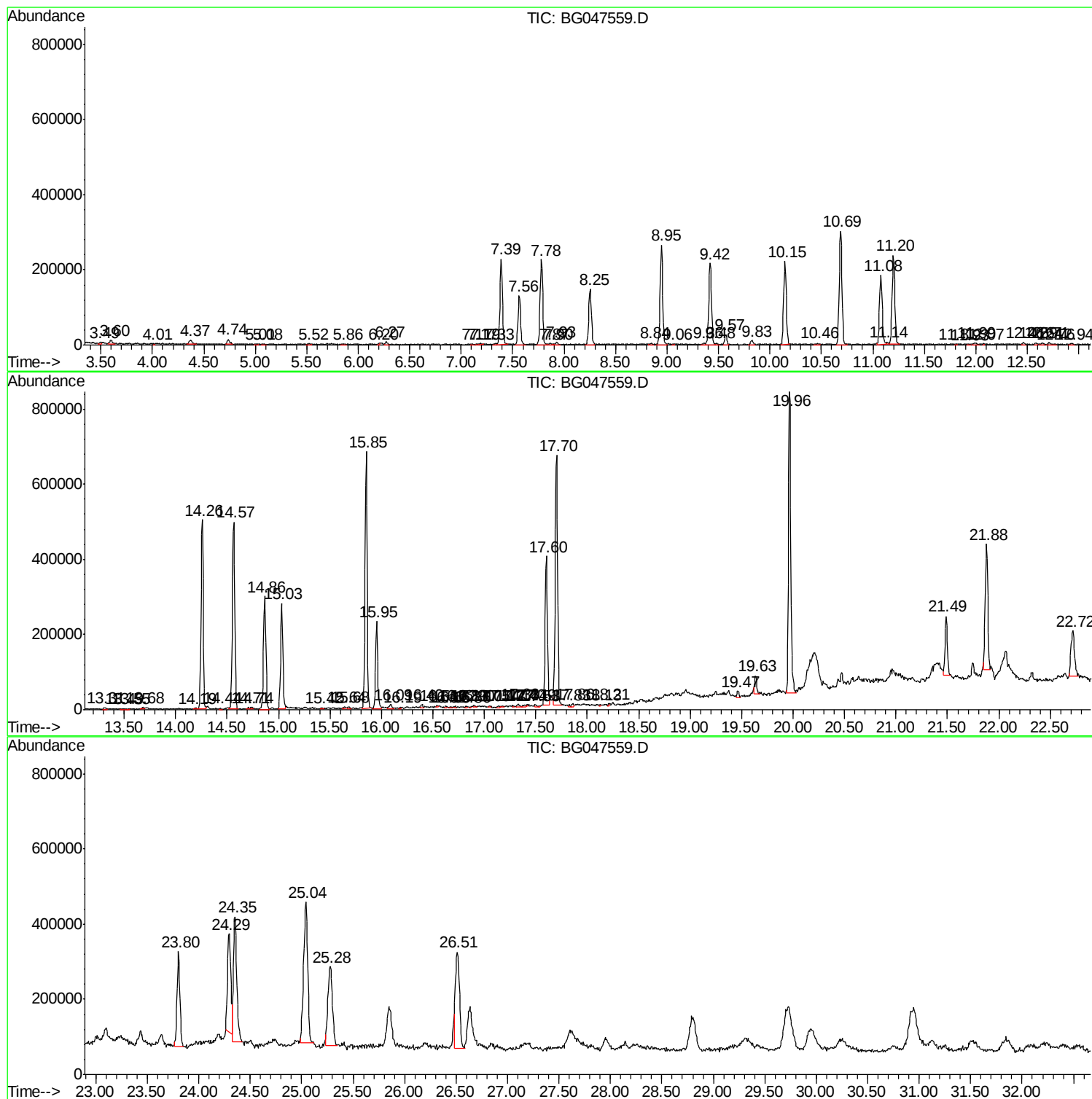
Sum of corrected areas: 16611829

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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



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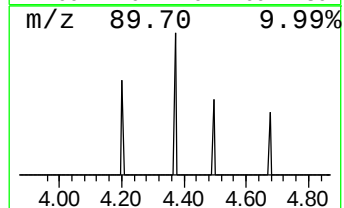
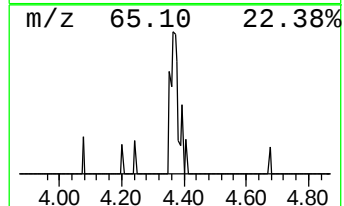
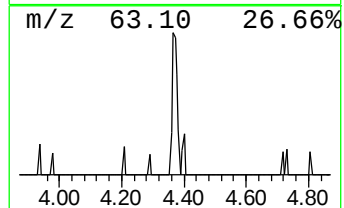
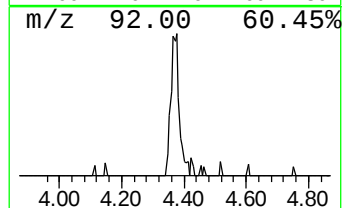
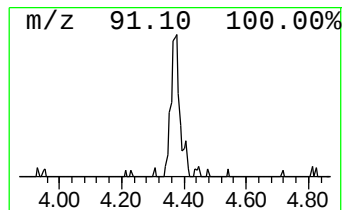
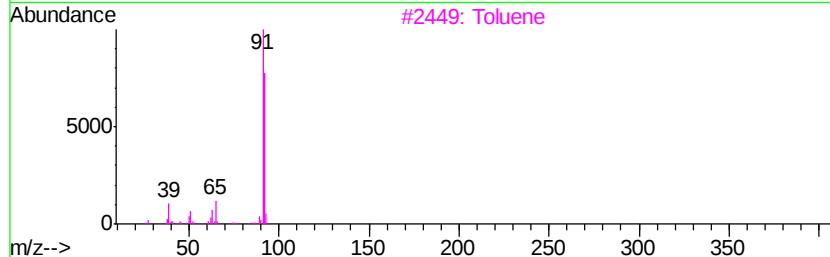
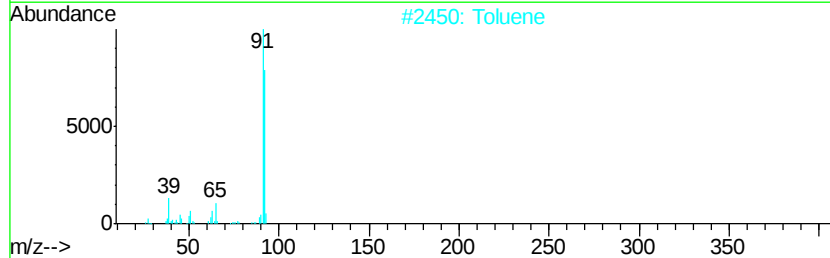
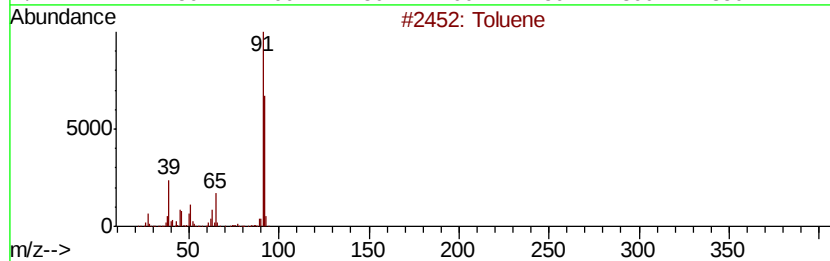
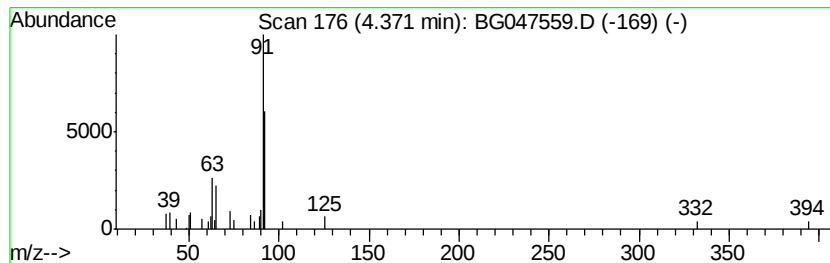
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	2.05 ng/ul	27433	1,4-Dichlorobenzene-d4	8.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Toluene	92	C7H8	000108-88-3	64
2		Toluene	92	C7H8	000108-88-3	64
3		Toluene	92	C7H8	000108-88-3	58
4		Toluene	92	C7H8	000108-88-3	58
5		Toluene	92	C7H8	000108-88-3	58



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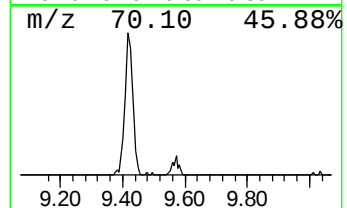
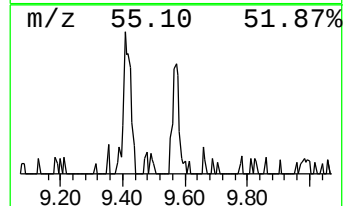
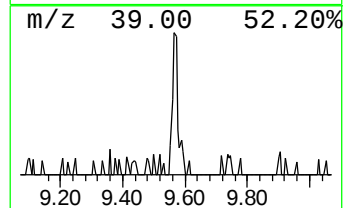
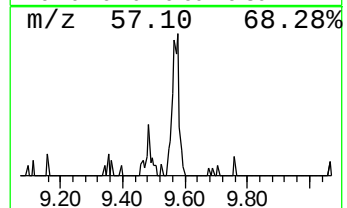
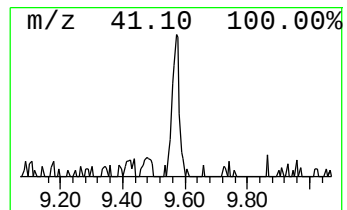
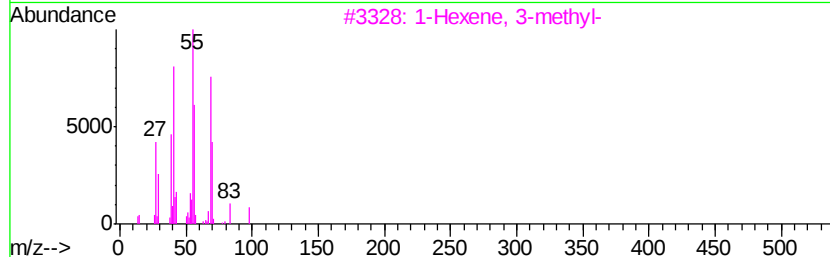
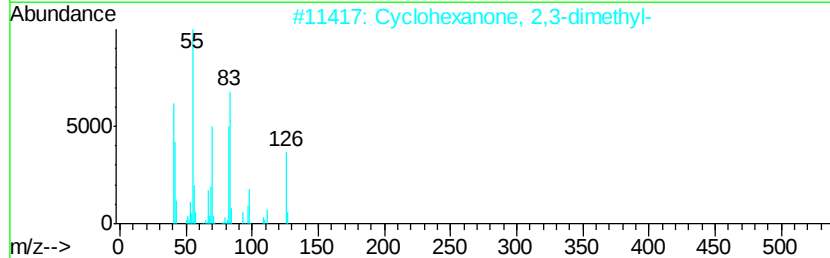
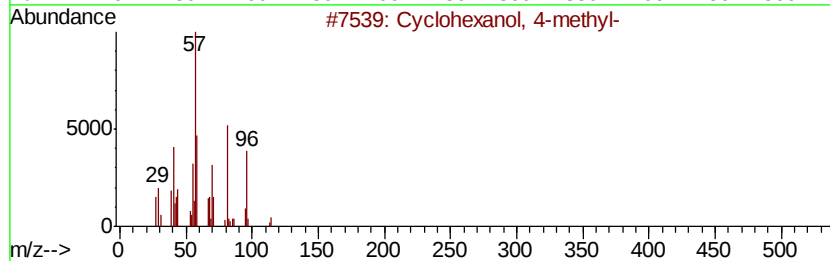
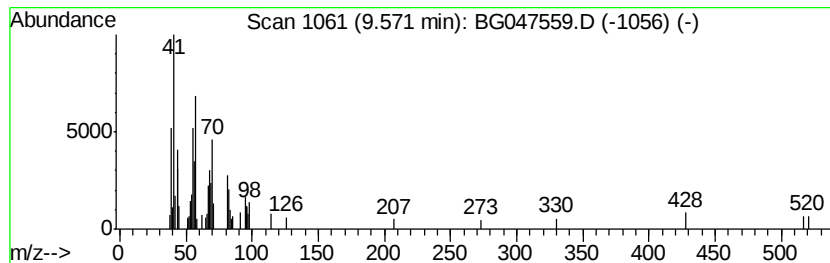
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.57	2.98 ng/ul	39869	1,4-Dichlorobenzene-d4	8.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanol, 4-methyl-	114	C7H14O	000589-91-3	43
2		Cyclohexanone, 2,3-dimethyl-	126	C8H14O	013395-76-1	35
3		1-Hexene, 3-methyl-	98	C7H14	003404-61-3	35
4		1-Hexene, 3-methyl-	98	C7H14	003404-61-3	35
5		2-Propen-1-amine, N-2-propenyl-	97	C6H11N	000124-02-7	30



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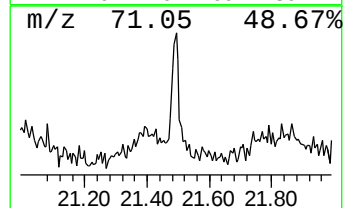
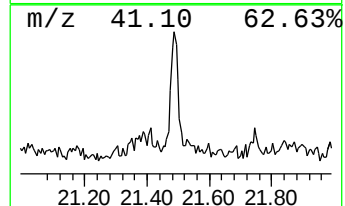
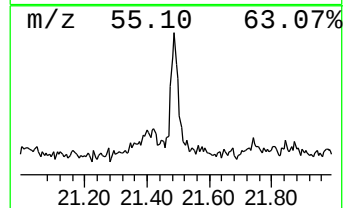
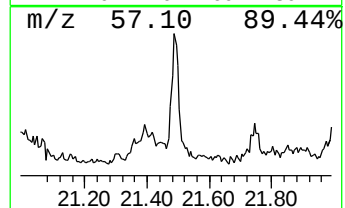
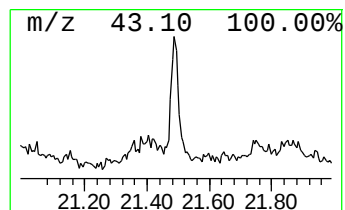
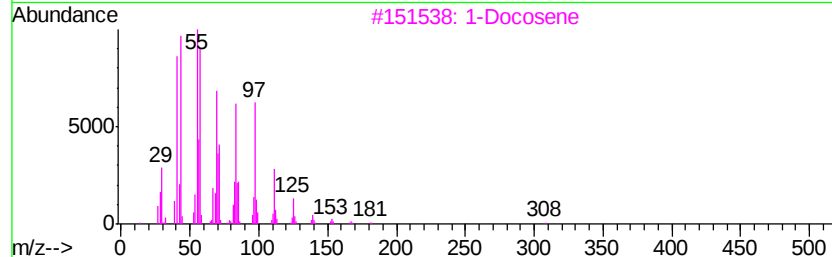
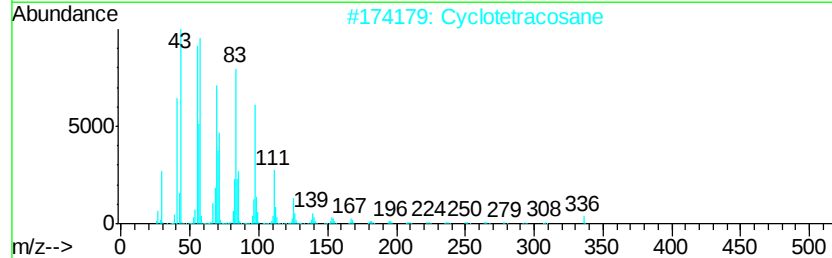
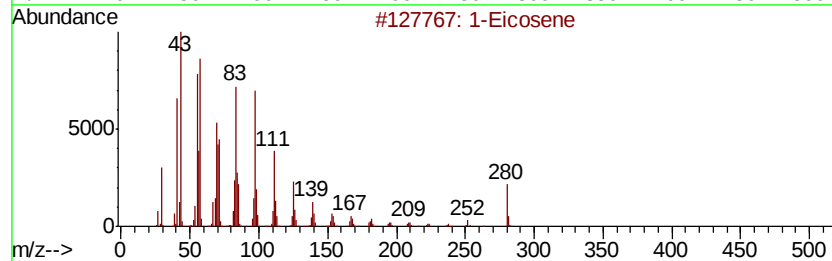
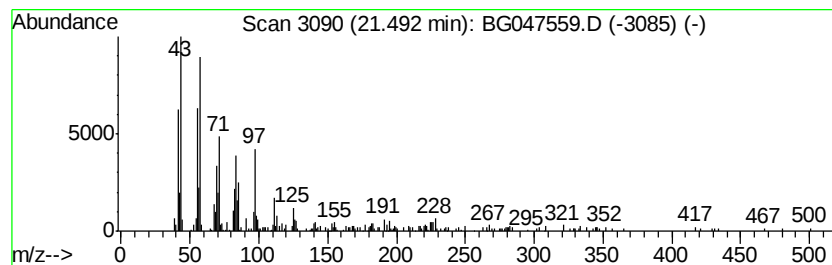
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TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.49	8.22 ng/ul	236112	Chrysene-d12	21.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Eicosene	280	C20H40	003452-07-1	89
2		Cyclotetracosane	336	C24H48	000297-03-0	83
3		1-Docosene	308	C22H44	001599-67-3	81
4		9-Tricosene, (Z)-	322	C23H46	027519-02-4	74
5		Cetene	224	C16H32	000629-73-2	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
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Acq On : 4 Dec 2020 5:32
Operator : CG/JU
Sample : L4855-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7Y9

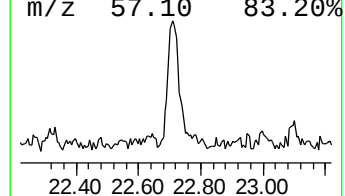
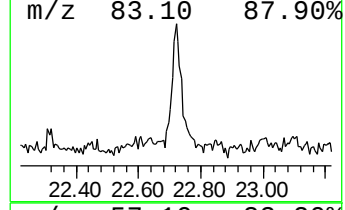
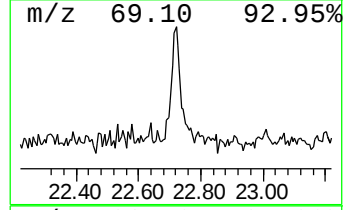
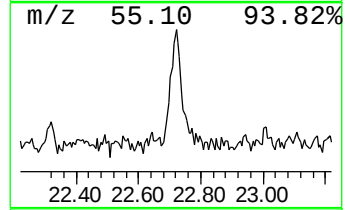
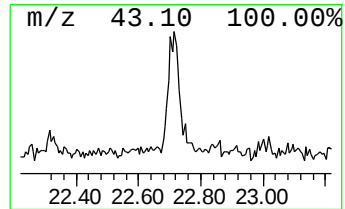
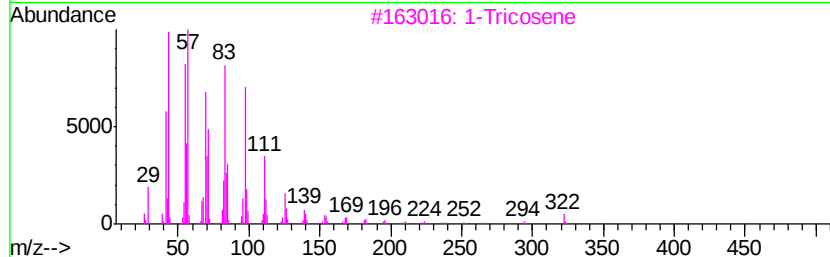
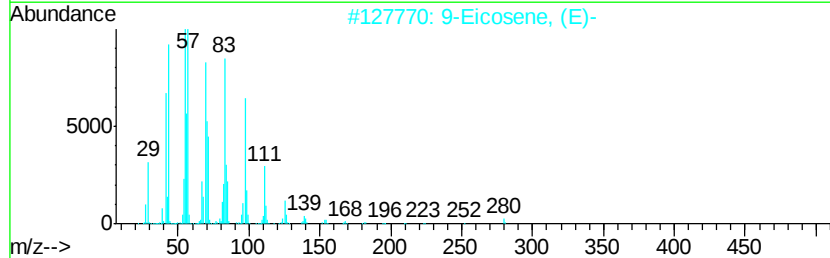
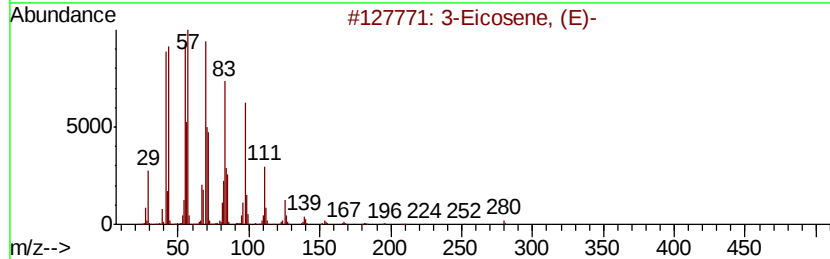
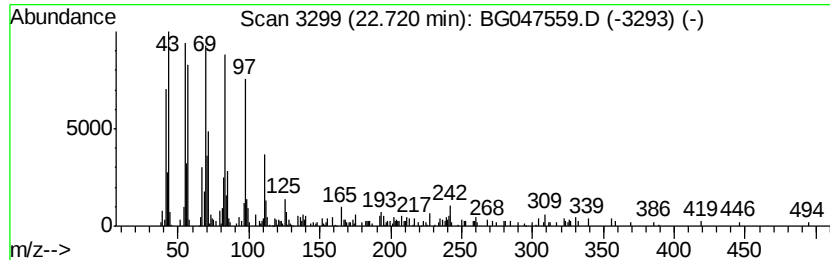
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.72	9.76 ng/ul	280351	Chrysene-d12	21.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2		9-Eicosene, (E)-	280	C20H40	074685-29-3	91
3		1-Tricosene	322	C23H46	018835-32-0	86
4		Cyclopropane, 1-(1,2-dimethylpro...	252	C18H36	041977-42-8	81
5		5-Eicosene, (E)-	280	C20H40	074685-30-6	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
Data File : BG047559.D
Acq On : 4 Dec 2020 5:32
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Sample : L4855-01
Misc :
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Instrument :
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ClientSampleId :
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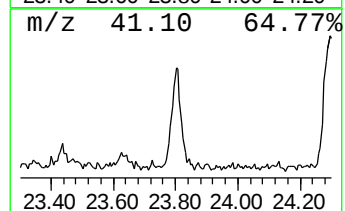
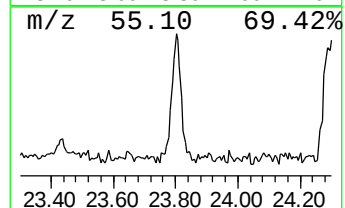
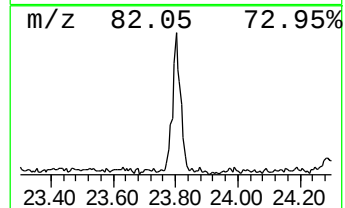
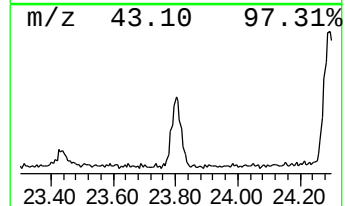
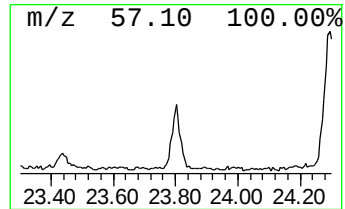
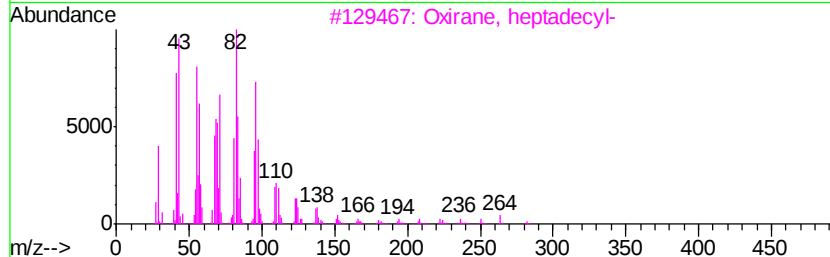
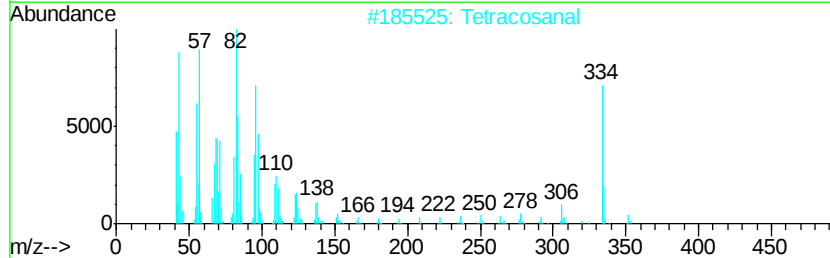
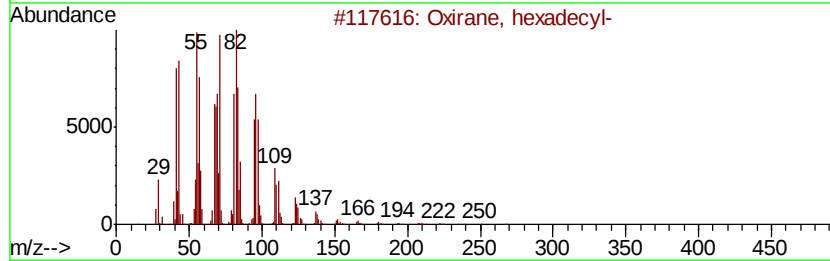
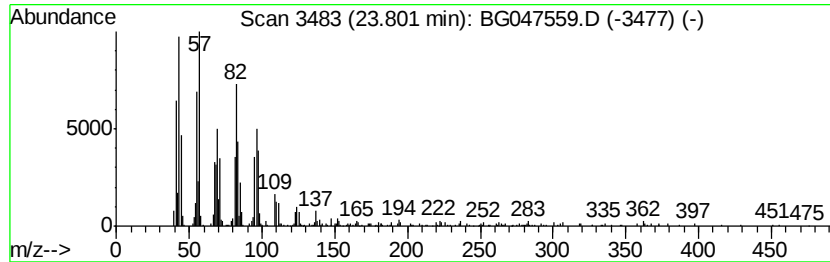
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Oxirane, hexadecyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.80	15.84 ng/ul	513705	Perylene-d12	25.28

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	74
2			Tetracosanal	352	C24H48O	057866-08-7	74
3			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	70
4			Octadecanal	268	C18H36O	000638-66-4	68
5			Pentadecanal-	226	C15H30O	002765-11-9	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
Data File : BG047559.D
Acq On : 4 Dec 2020 5:32
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Sample : L4855-01
Misc :
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Instrument :
BNA_G
ClientSampleId :
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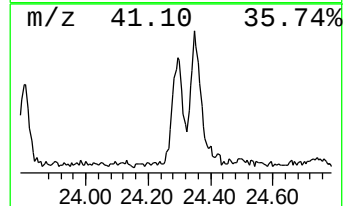
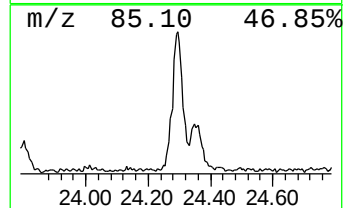
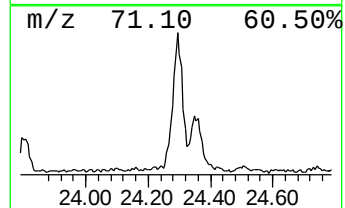
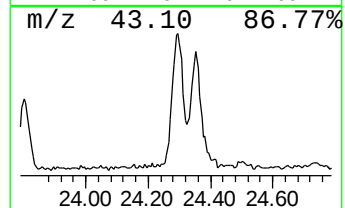
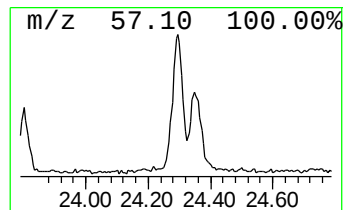
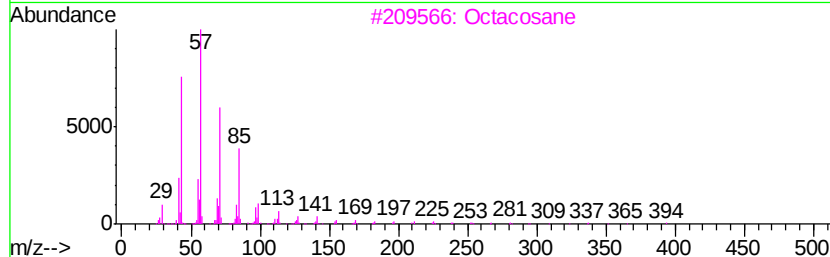
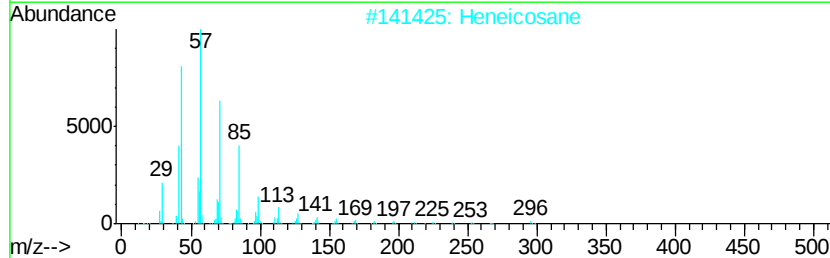
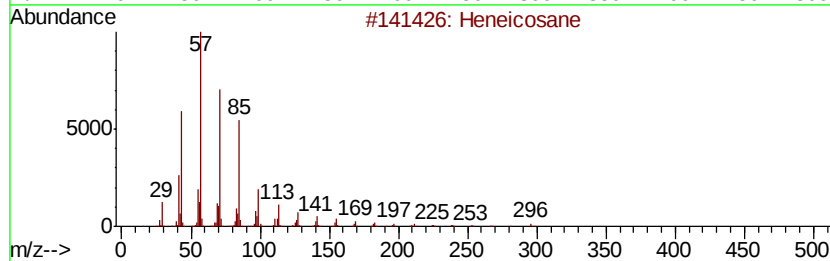
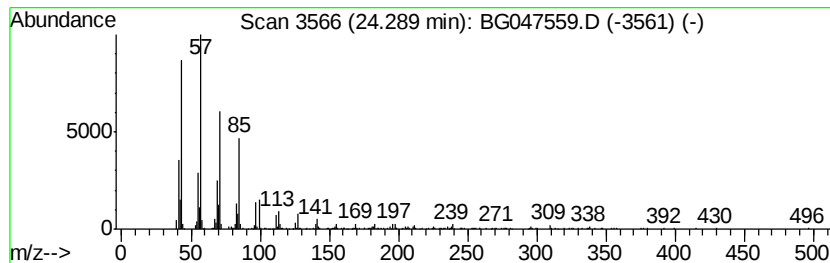
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.29	18.20 ng/ul	589991	Perylene-d12	25.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	93
2		Heneicosane	296	C21H44	000629-94-7	93
3		Octacosane	394	C28H58	000630-02-4	90
4		Eicosane	282	C20H42	000112-95-8	90
5		Tricosane	324	C23H48	000638-67-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
Data File : BG047559.D
Acq On : 4 Dec 2020 5:32
Operator : CG/JU
Sample : L4855-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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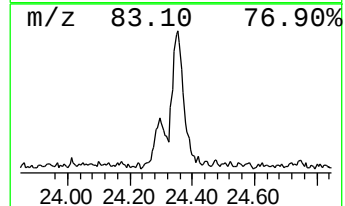
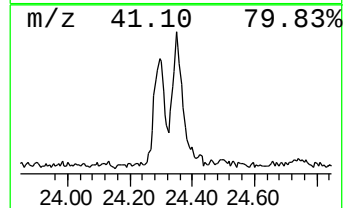
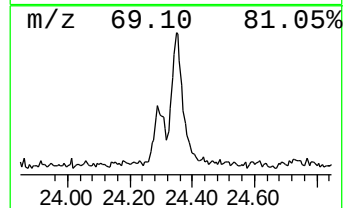
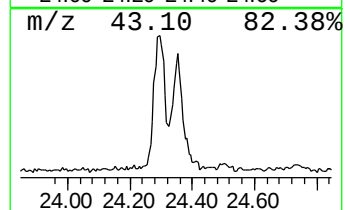
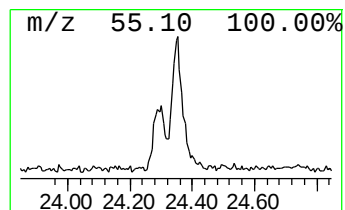
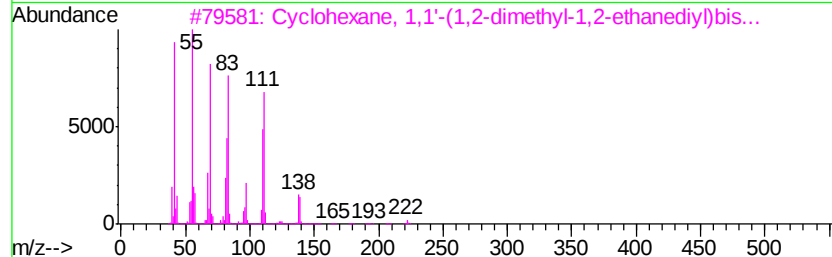
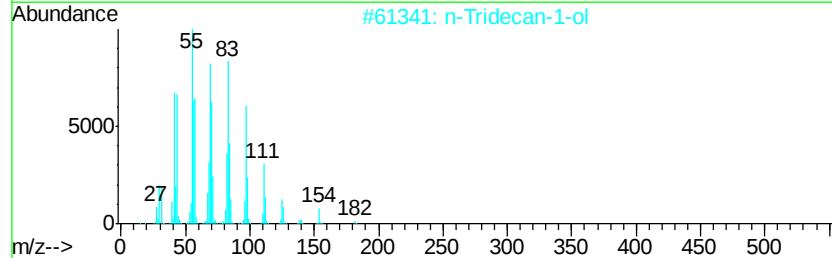
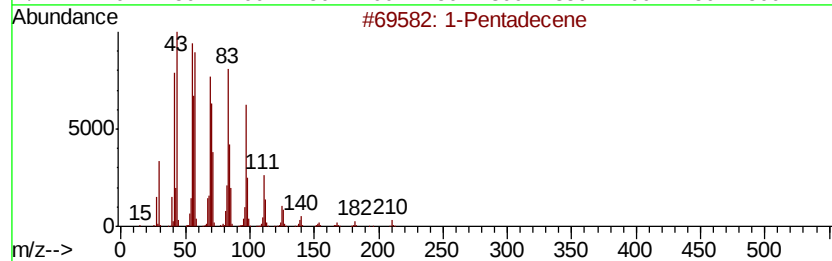
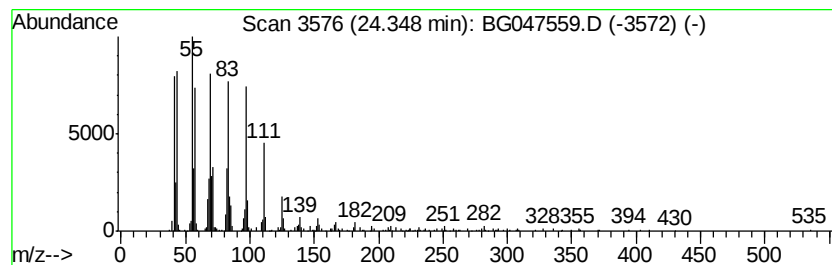
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.35	23.75 ng/ul	770236	Perylene-d12	25.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentadecene	210	C15H30	013360-61-7	81
2		n-Tridecan-1-ol	200	C13H28O	000112-70-9	80
3		Cyclohexane, 1,1'-(1,2-dimethyl-...	222	C16H30	054889-87-1	76
4		Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	70
5		9-Tricosene, (Z)-	322	C23H46	027519-02-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\
Data File : BG047559.D
Acq On : 4 Dec 2020 5:32
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Sample : L4855-01
Misc :
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ClientSampleId :
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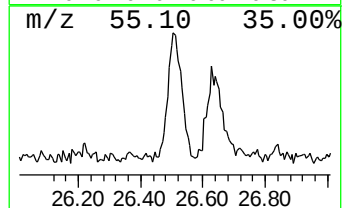
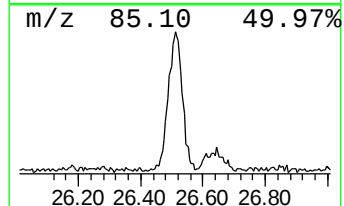
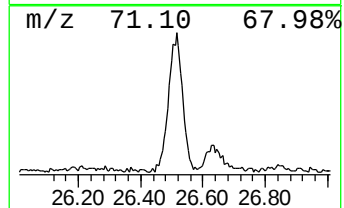
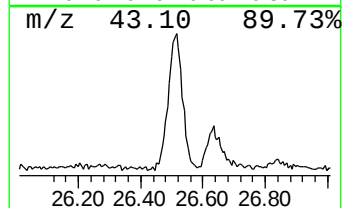
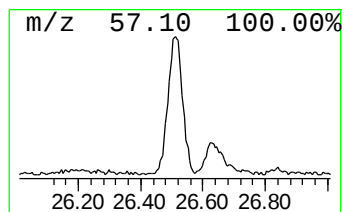
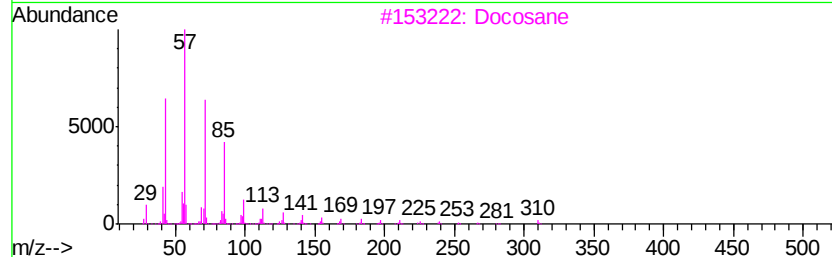
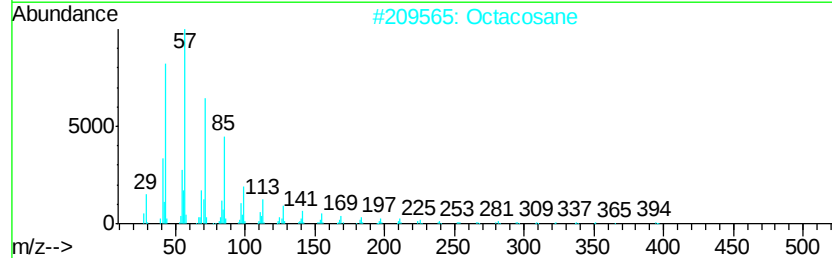
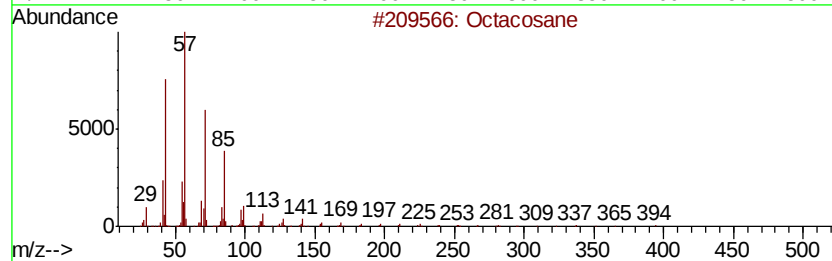
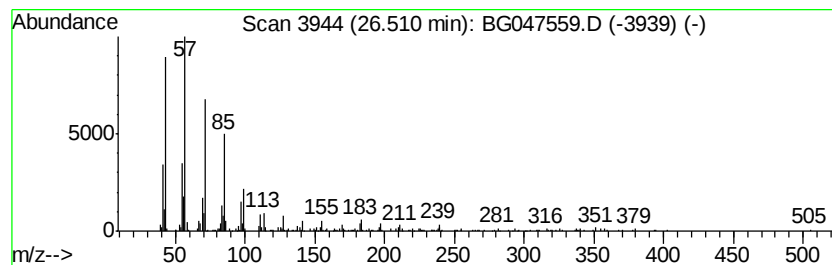
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.51	21.99 ng/ul	713168	Perylene-d12	25.28

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	96
2		Octacosane	394	C28H58	000630-02-4	95
3		Docosane	310	C22H46	000629-97-0	94
4		Tetratetracontane	619	C44H90	007098-22-8	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG120320\
Data File : BG047559.D
Acq On : 4 Dec 2020 5:32
Operator : CG/JU
Sample : L4855-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	9.57	3.0	ng/ul	39869	1	8.25	268025	20.0
(DEL) Alkane: Str...	21.49	8.2	ng/ul	236112	5	21.88	574660	20.0
(DEL) Alkane: Str...	22.72	9.8	ng/ul	280351	5	21.88	574660	20.0
Oxirane, hexadecyl-	23.80	15.8	ng/ul	513705	6	25.28	648495	20.0
(DEL) Alkane: Str...	24.29	18.2	ng/ul	589991	6	25.28	648495	20.0
(DEL) Alkane: Str...	24.35	23.8	ng/ul	770236	6	25.28	648495	20.0
(DEL) Alkane: Str...	26.51	22.0	ng/ul	713168	6	25.28	648495	20.0