

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072621\
 Data File : BG049499.D
 Acq On : 27 Jul 2021 9:00
 Operator : CG/JU
 Sample : PB137867BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK867

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-BG072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BG049499.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.079	3	7	16	rBV4	19638	39669	5.36%	0.505%
2	3.162	18	21	26	rBV6	3817	5328	0.72%	0.068%
3	3.432	63	67	74	rBV3	6709	9800	1.32%	0.125%
4	3.614	97	98	101	rVB2	1204	1019	0.14%	0.013%
5	3.690	107	111	113	rVB3	1542	1471	0.20%	0.019%
6	3.861	137	140	151	rBV	61107	113378	15.31%	1.445%
7	4.084	176	178	182	rBV3	1963	2341	0.32%	0.030%
8	4.195	194	197	203	rVB3	1811	2467	0.33%	0.031%
9	4.454	239	241	243	rVB	1455	1099	0.15%	0.014%
10	4.771	291	295	299	rBV3	1164	2067	0.28%	0.026%
11	4.865	308	311	313	rVB2	847	1047	0.14%	0.013%
12	4.942	321	324	327	rVB3	1038	1280	0.17%	0.016%
13	5.364	391	396	397	rBV3	1187	1964	0.27%	0.025%
14	5.394	400	401	404	rBV2	1392	1589	0.21%	0.020%
15	5.500	416	419	422	rVB3	1129	1369	0.18%	0.017%
16	5.570	430	431	434	rVB	1553	1253	0.17%	0.016%
17	5.599	434	436	438	rBV	989	1178	0.16%	0.015%
18	5.805	468	471	474	rVB2	782	1020	0.14%	0.013%
19	5.928	489	492	495	rVB	986	1365	0.18%	0.017%
20	6.351	562	564	566	rVB	1221	1064	0.14%	0.014%
21	6.410	569	574	576	rBV	954	1171	0.16%	0.015%
22	6.774	633	636	637	rBV	1358	1210	0.16%	0.015%
23	7.215	704	711	719	rBV	112638	199803	26.98%	2.546%
24	7.274	719	721	724	rVB	1477	1103	0.15%	0.014%
25	7.379	732	739	748	rBV	70977	124098	16.76%	1.581%
26	7.585	766	774	783	rVB3	108231	191467	25.85%	2.440%
27	8.055	848	854	863	rVB	87467	152942	20.65%	1.949%
28	8.196	874	878	879	rVB	898	1167	0.16%	0.015%
29	8.766	967	975	985	rBV	136968	240392	32.46%	3.063%
30	8.965	1006	1009	1015	rVB	875	1435	0.19%	0.018%
31	9.230	1047	1054	1063	rVB	105055	195399	26.39%	2.490%
32	9.465	1091	1094	1097	rBV	963	1075	0.15%	0.014%
33	9.952	1170	1177	1185	rVB2	94918	177751	24.00%	2.265%
34	10.499	1261	1270	1277	rBV2	133680	247474	33.42%	3.153%
35	10.552	1277	1279	1282	rVB2	1599	1529	0.21%	0.019%
36	10.881	1328	1335	1343	rVB	119213	217847	29.42%	2.776%

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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-BG072421.M
 Title : SVOA CALIBRATION

37	11.010	1351	1357	1365	rBV	122403	235114	31.75%	2.996%
38	11.973	1520	1521	1526	rBV	756	1033	0.14%	0.013%
39	12.384	1589	1591	1596	rBV	925	1189	0.16%	0.015%
40	12.472	1601	1606	1607	rBV2	1632	2668	0.36%	0.034%

41	13.201	1725	1730	1731	rBV2	1230	1342	0.18%	0.017%
42	14.106	1877	1884	1892	rBV	258853	384541	51.93%	4.900%
43	14.399	1927	1934	1941	rBV	278727	446048	60.23%	5.683%
44	14.705	1979	1986	1994	rBV	202245	321924	43.47%	4.102%
45	14.899	2013	2019	2028	rBV	132789	228253	30.82%	2.908%

46	15.069	2046	2048	2051	rVB	1403	1491	0.20%	0.019%
47	15.098	2051	2053	2060	rBV2	1053	2070	0.28%	0.026%
48	15.698	2149	2155	2163	rBV	389833	580397	78.37%	7.395%
49	15.815	2168	2175	2187	rVV	129190	192031	25.93%	2.447%
50	15.891	2187	2188	2193	rVB2	1598	1142	0.15%	0.015%

51	15.938	2193	2196	2202	rBV2	4066	5507	0.74%	0.070%
52	16.285	2252	2255	2258	rVB2	819	1212	0.16%	0.015%
53	17.237	2416	2417	2424	rVB	1120	1606	0.22%	0.020%
54	17.454	2448	2454	2461	rVB	274433	428007	57.79%	5.453%
55	17.554	2464	2471	2479	rVV	445838	683656	92.32%	8.711%

56	17.718	2496	2499	2502	rBV	1281	1875	0.25%	0.024%
57	17.830	2514	2518	2520	rBV	1028	1207	0.16%	0.015%
58	18.353	2605	2607	2610	rVB	890	1117	0.15%	0.014%
59	18.881	2696	2697	2702	rBV	1021	1202	0.16%	0.015%
60	19.081	2728	2731	2737	rVB3	1163	1796	0.24%	0.023%

61	19.134	2737	2740	2744	rBV3	1080	1432	0.19%	0.018%
62	19.404	2782	2786	2790	rBV4	5409	7356	0.99%	0.094%
63	19.475	2794	2798	2802	rVV3	6422	7425	1.00%	0.095%
64	19.534	2805	2808	2810	rBV2	788	1057	0.14%	0.013%
65	19.692	2832	2835	2837	rBV2	1369	1264	0.17%	0.016%

66	19.722	2837	2840	2842	rVV	1040	1268	0.17%	0.016%
67	19.757	2842	2846	2849	rVV4	1409	1991	0.27%	0.025%
68	19.839	2854	2860	2867	rBV	528941	740561	100.00%	9.436%
69	19.951	2876	2879	2881	rBV	1183	1198	0.16%	0.015%
70	20.009	2888	2889	2892	rBV2	1239	1171	0.16%	0.015%

71	20.045	2894	2895	2897	rBV	2298	1956	0.26%	0.025%
72	20.074	2899	2900	2902	rBV2	2821	2152	0.29%	0.027%
73	20.168	2915	2916	2917	rBV	2700	1520	0.21%	0.019%
74	20.186	2917	2919	2921	rVV	3050	2599	0.35%	0.033%
75	20.315	2940	2941	2942	rBV	2667	1115	0.15%	0.014%

76	20.332	2942	2944	2946	rVV2	2204	1883	0.25%	0.024%
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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

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
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77	20.421	2957	2959	2961	rBV2	2427	2398	0.32%	0.031%
78	20.526	2975	2977	2978	rVB2	2492	1608	0.22%	0.020%
79	20.544	2978	2980	2981	rBV	1869	1178	0.16%	0.015%
80	20.714	3007	3009	3010	rBV2	2256	1495	0.20%	0.019%
81	21.266	3102	3103	3105	rBV2	3398	2407	0.33%	0.031%
82	21.736	3175	3183	3189	rBV	252741	431808	58.31%	5.502%
83	24.227	3605	3607	3611	rBV3	4072	5214	0.70%	0.066%
84	24.791	3693	3703	3713	rBV2	261846	695690	93.94%	8.864%
85	25.020	3733	3742	3753	rVB2	149326	422213	57.01%	5.380%
86	25.220	3774	3776	3779	rBV4	3110	2396	0.32%	0.031%
87	25.831	3878	3880	3885	rVB4	3444	4856	0.66%	0.062%
88	26.606	4011	4012	4013	rVB	3539	1247	0.17%	0.016%
89	26.718	4029	4031	4032	rVB2	3197	1397	0.19%	0.018%
90	28.145	4272	4274	4276	rVB3	3595	2308	0.31%	0.029%
91	28.169	4276	4278	4279	rBV	3325	2349	0.32%	0.030%
92	28.944	4408	4410	4414	rBV4	4082	3548	0.48%	0.045%
93	29.185	4449	4451	4452	rBV2	3333	2158	0.29%	0.027%
94	29.532	4508	4510	4511	rBV2	3547	1658	0.22%	0.021%
95	30.066	4599	4601	4602	rVB2	3864	1910	0.26%	0.024%
96	30.748	4715	4717	4718	rBV2	3291	1403	0.19%	0.018%
97	31.235	4798	4800	4801	rBV2	3433	2357	0.32%	0.030%
98	31.253	4801	4803	4804	rVV	4309	2557	0.35%	0.033%
99	32.099	4945	4947	4949	rBV3	3492	2164	0.29%	0.028%
100	32.304	4980	4982	4983	rBV2	3424	2188	0.30%	0.028%

Sum of corrected areas: 7848514

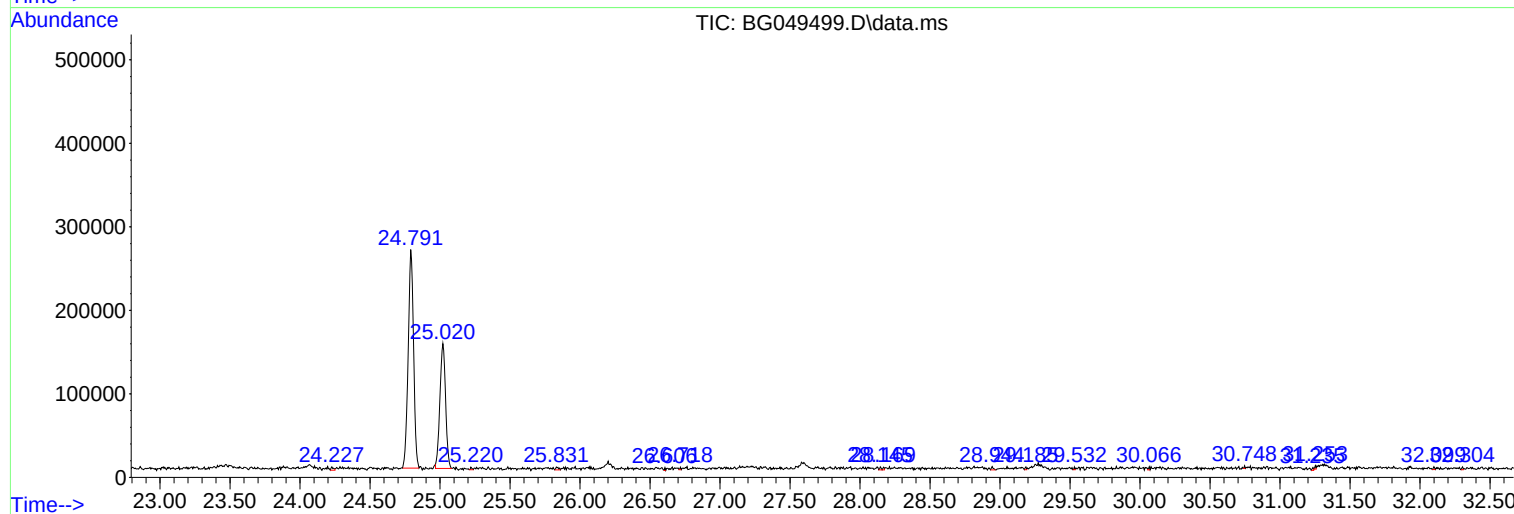
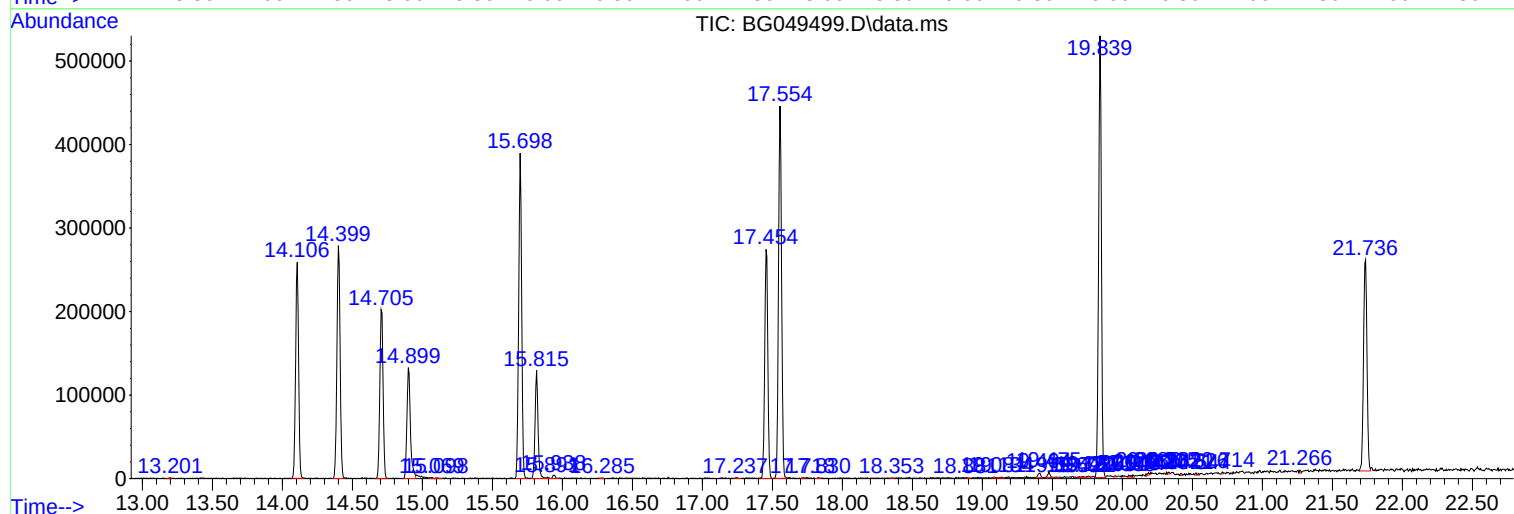
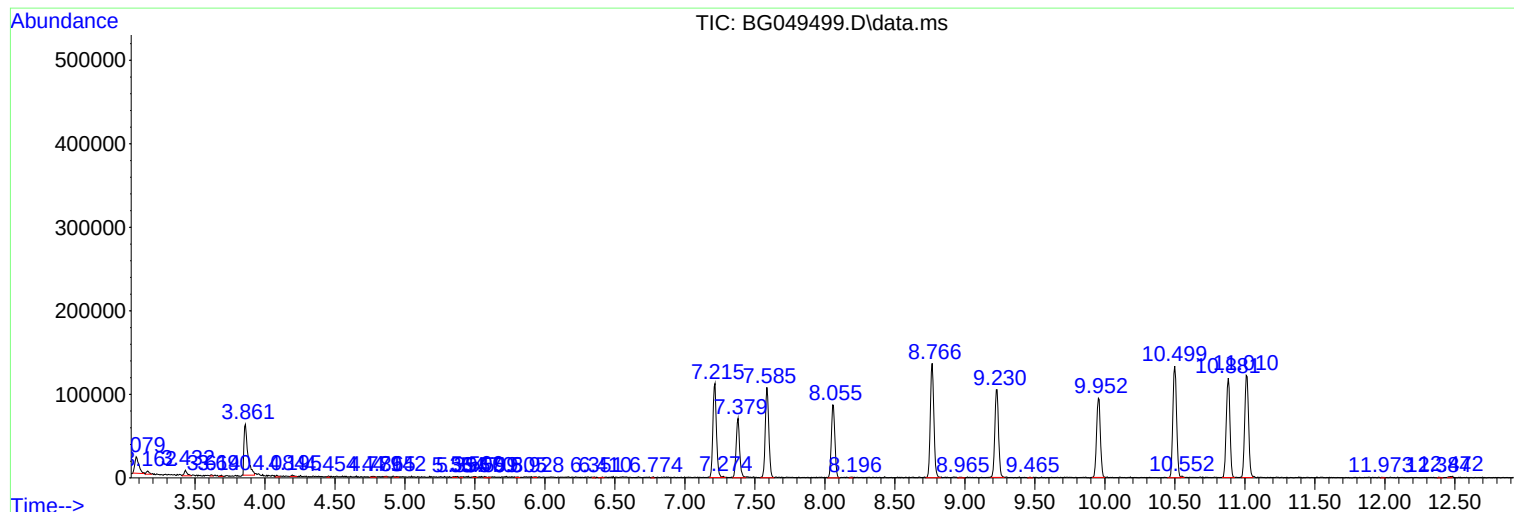
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P



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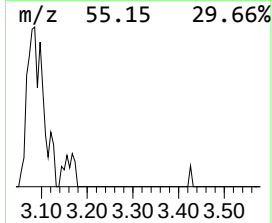
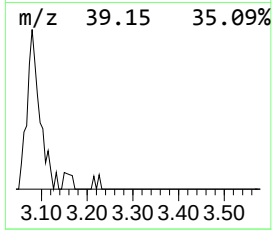
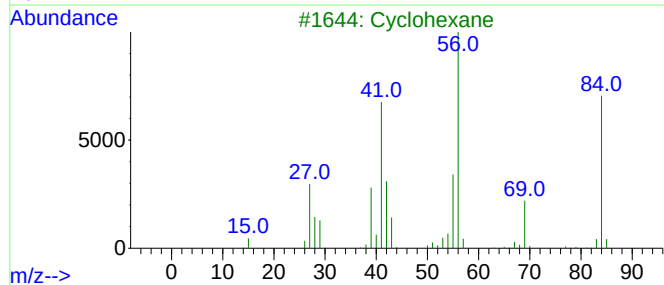
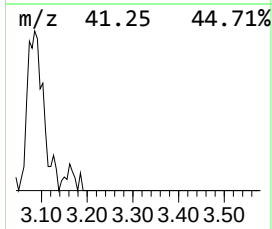
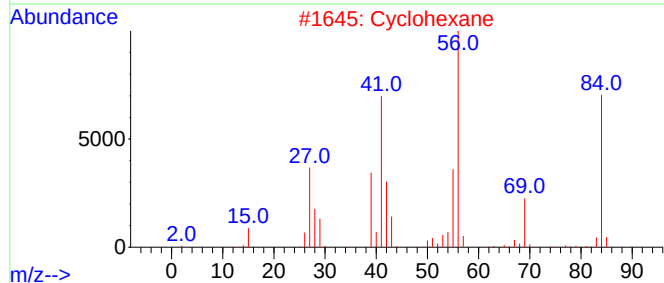
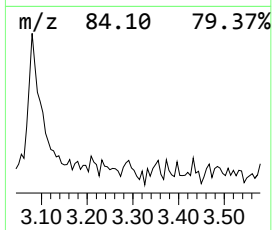
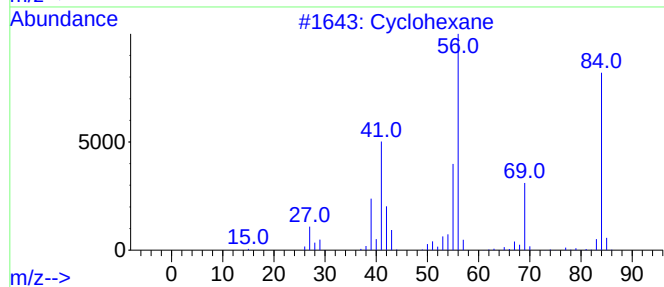
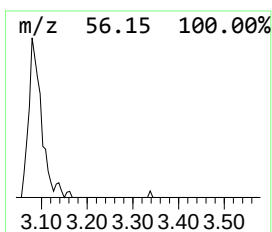
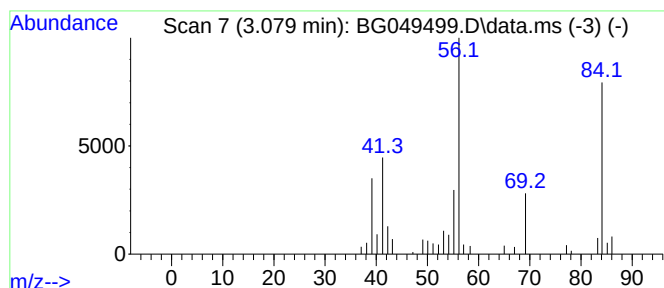
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.079 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.079	5.19 ng/ul	39669	1,4-Dichlorobenzene-d4	8.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane	84	C6H12	000110-82-7	87
2			Cyclohexane	84	C6H12	000110-82-7	80
3			Cyclohexane	84	C6H12	000110-82-7	72
4			Cyclohexane	84	C6H12	000110-82-7	72
5			2-Amino-oxazole	84	C3H4N2O	004570-45-0	72



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	3.079	5.2	ng/u1	39669	1	8.055	152942	20.0