

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072921\
 Data File : BG049553.D
 Acq On : 29 Jul 2021 10:27
 Operator : CG/JU
 Sample : PB137954BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK954

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: BG049553.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.085	3	8	15	rVB5	19644	41793	5.24%	0.503%
2	3.432	63	67	73	rVB4	6472	11225	1.41%	0.135%
3	3.784	125	127	130	rVB2	1559	1577	0.20%	0.019%
4	3.855	136	139	158	rBV	77331	158281	19.85%	1.904%
5	4.495	246	248	255	rVB4	1748	3005	0.38%	0.036%
6	4.712	282	285	287	rBV2	1117	1706	0.21%	0.021%
7	5.165	360	362	366	rBV2	1569	1282	0.16%	0.015%
8	5.353	393	394	397	rBV3	1040	1406	0.18%	0.017%
9	5.594	433	435	440	rVB2	1174	1202	0.15%	0.014%
10	5.693	447	452	453	rBV	1217	1521	0.19%	0.018%
11	6.146	527	529	531	rBV	1162	1154	0.14%	0.014%
12	6.669	616	618	622	rVB	1393	1356	0.17%	0.016%
13	6.815	638	643	644	rBV2	1004	1125	0.14%	0.014%
14	6.833	644	646	650	rBV	758	1175	0.15%	0.014%
15	7.203	701	709	718	rBV2	131125	229905	28.83%	2.766%
16	7.291	721	724	727	rVB2	1376	1238	0.16%	0.015%
17	7.368	731	737	750	rVB	79202	142483	17.87%	1.714%
18	7.579	766	773	780	rBV2	123446	221690	27.80%	2.667%
19	8.049	846	853	862	rVB	81347	137926	17.30%	1.659%
20	8.648	951	955	956	rBV2	1022	1158	0.15%	0.014%
21	8.754	967	973	982	rVB2	155717	273404	34.28%	3.289%
22	9.218	1045	1052	1061	rBV	130073	233730	29.31%	2.812%
23	9.318	1065	1069	1072	rBV3	907	1333	0.17%	0.016%
24	9.424	1080	1087	1088	rBV2	1446	1631	0.20%	0.020%
25	9.946	1168	1176	1186	rBV2	116520	216540	27.15%	2.605%
26	10.487	1262	1268	1277	rBV	155820	289328	36.28%	3.481%
27	10.875	1326	1334	1341	rBV	111253	197440	24.76%	2.375%
28	10.922	1341	1342	1345	rVB	1674	1119	0.14%	0.013%
29	11.004	1345	1356	1368	rBV3	148257	285200	35.76%	3.431%
30	11.697	1472	1474	1477	rBV3	926	1152	0.14%	0.014%
31	11.967	1518	1520	1523	rBV	979	1115	0.14%	0.013%
32	12.455	1599	1603	1604	rBV2	1864	1804	0.23%	0.022%
33	12.596	1625	1627	1630	rVB2	1111	1198	0.15%	0.014%
34	13.077	1703	1709	1712	rBV2	823	1609	0.20%	0.019%
35	13.659	1805	1808	1811	rBV2	1023	1285	0.16%	0.015%
36	14.100	1875	1883	1892	rBV	298274	453977	56.93%	5.462%

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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

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	14.393	1927	1933	1941	rVB	311469	497615	62.40%	5.987%
38	14.699	1979	1985	1992	rVB	183661	289607	36.32%	3.484%
39	14.893	2012	2018	2030	rBV	134487	242429	30.40%	2.917%
40	15.139	2057	2060	2063	rVB	1281	1096	0.14%	0.013%
41	15.692	2147	2154	2161	rBV	412805	634356	79.55%	7.632%
42	15.809	2168	2174	2184	rBV2	127489	204322	25.62%	2.458%
43	15.938	2191	2196	2199	rVB3	4717	6992	0.88%	0.084%
44	16.743	2330	2333	2335	rBV3	1026	1184	0.15%	0.014%
45	17.054	2385	2386	2391	rBV3	1223	1773	0.22%	0.021%
46	17.231	2414	2416	2421	rVB	1270	1543	0.19%	0.019%
47	17.448	2446	2453	2460	rBV	235062	360343	45.19%	4.335%
48	17.548	2463	2470	2478	rBV	473017	725136	90.93%	8.724%
49	17.707	2494	2497	2502	rBV	709	1163	0.15%	0.014%
50	17.842	2515	2520	2522	rBV2	994	1388	0.17%	0.017%
51	18.494	2626	2631	2635	rVB	951	1751	0.22%	0.021%
52	18.594	2640	2648	2653	rVB2	806	1894	0.24%	0.023%
53	18.641	2653	2656	2659	rBV2	920	1110	0.14%	0.013%
54	18.799	2681	2683	2688	rVB	1233	1338	0.17%	0.016%
55	19.064	2726	2728	2731	rBV	1177	1176	0.15%	0.014%
56	19.093	2731	2733	2736	rVV	860	1216	0.15%	0.015%
57	19.128	2736	2739	2742	rVB2	1018	1246	0.16%	0.015%
58	19.404	2778	2786	2789	rBV4	5459	9343	1.17%	0.112%
59	19.469	2794	2797	2801	rBV2	5935	6914	0.87%	0.083%
60	19.839	2853	2860	2867	rBV	563337	797455	100.00%	9.594%
61	19.992	2882	2886	2889	rVB2	1390	1638	0.21%	0.020%
62	20.092	2900	2903	2905	rBV3	1517	1755	0.22%	0.021%
63	20.174	2915	2917	2918	rBV	2070	1276	0.16%	0.015%
64	20.274	2932	2934	2936	rVB2	2593	1969	0.25%	0.024%
65	20.426	2958	2960	2961	rVB2	2478	1509	0.19%	0.018%
66	20.450	2961	2964	2967	rBV2	1789	2672	0.34%	0.032%
67	20.585	2985	2987	2988	rVB2	2538	1091	0.14%	0.013%
68	20.614	2990	2992	2994	rVB2	1845	1176	0.15%	0.014%
69	20.644	2994	2997	2999	rBV2	1881	2356	0.30%	0.028%
70	20.661	2999	3000	3004	rVV3	2483	2258	0.28%	0.027%
71	20.708	3006	3008	3011	rVB2	1898	1430	0.18%	0.017%
72	20.814	3024	3026	3028	rVB3	3390	2440	0.31%	0.029%
73	20.838	3028	3030	3031	rBV	2286	2099	0.26%	0.025%
74	21.466	3135	3137	3141	rBV4	3152	3893	0.49%	0.047%
75	21.543	3148	3150	3151	rBV2	2379	1603	0.20%	0.019%
76	21.730	3176	3182	3189	rBV2	221090	367652	46.10%	4.423%

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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-BG072421.M
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77	21.848	3200	3202	3205	rBV4	2540	2278	0.29%	0.027%
78	22.829	3367	3369	3370	rVB2	3432	2020	0.25%	0.024%
79	24.110	3585	3587	3588	rBV	2350	1128	0.14%	0.014%
80	24.685	3684	3685	3686	rBV	3762	1696	0.21%	0.020%
81	24.791	3692	3703	3714	rVB	272029	754274	94.59%	9.075%
82	25.014	3731	3741	3752	rVB	131020	377247	47.31%	4.539%
83	25.337	3795	3796	3798	rBV2	2455	1647	0.21%	0.020%
84	25.449	3813	3815	3816	rBV2	2733	1738	0.22%	0.021%
85	26.183	3936	3940	3946	rBV7	7038	16168	2.03%	0.195%
86	26.671	4020	4023	4024	rBV3	2946	3223	0.40%	0.039%
87	28.380	4312	4314	4317	rVB3	3647	2949	0.37%	0.035%
88	28.539	4339	4341	4345	rBV4	2642	3964	0.50%	0.048%
89	28.586	4347	4349	4350	rBV2	3032	1644	0.21%	0.020%
90	29.162	4446	4447	4449	rBV2	2926	2439	0.31%	0.029%
91	29.925	4575	4577	4579	rBV2	3052	2543	0.32%	0.031%
92	30.495	4672	4674	4675	rBV2	2724	2245	0.28%	0.027%
93	30.971	4753	4755	4757	rBV3	3004	2834	0.36%	0.034%
94	31.147	4783	4785	4786	rBV2	2350	1605	0.20%	0.019%
95	31.312	4811	4813	4814	rBV2	2920	1267	0.16%	0.015%
96	31.441	4833	4835	4838	rBV4	4043	5141	0.64%	0.062%
97	31.817	4897	4899	4900	rBV	2973	1696	0.21%	0.020%
98	31.946	4919	4921	4923	rBV3	2508	2256	0.28%	0.027%
99	32.105	4946	4948	4949	rBV2	2290	1742	0.22%	0.021%
100	32.375	4992	4994	4998	rBV4	3640	4773	0.60%	0.057%

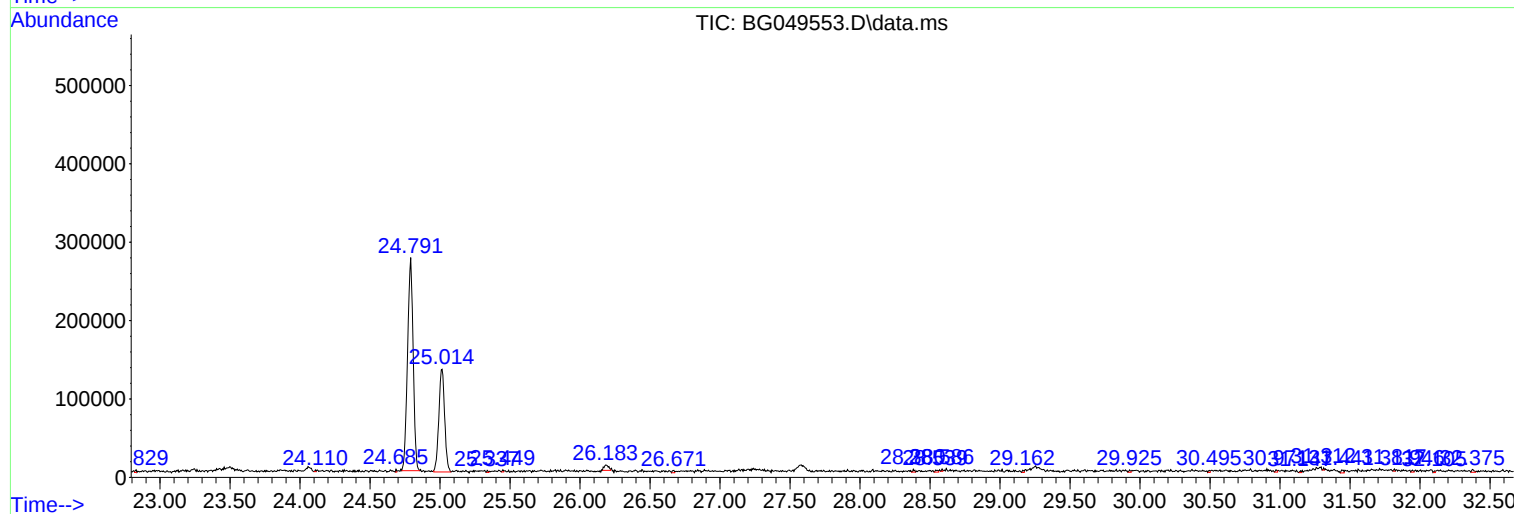
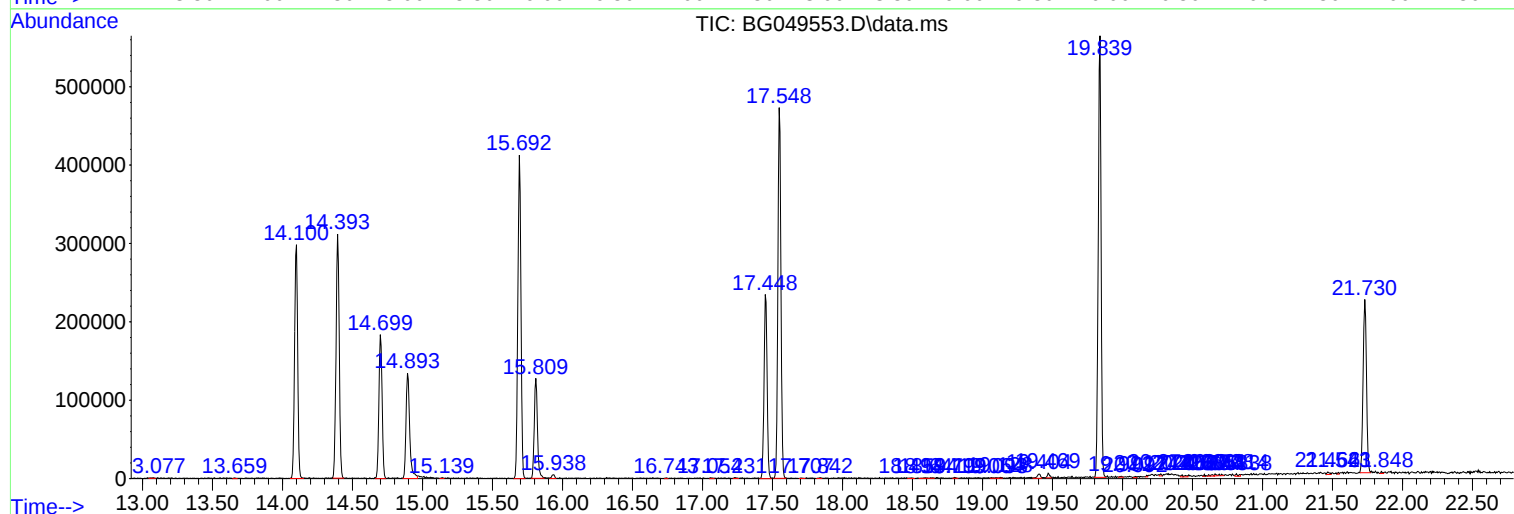
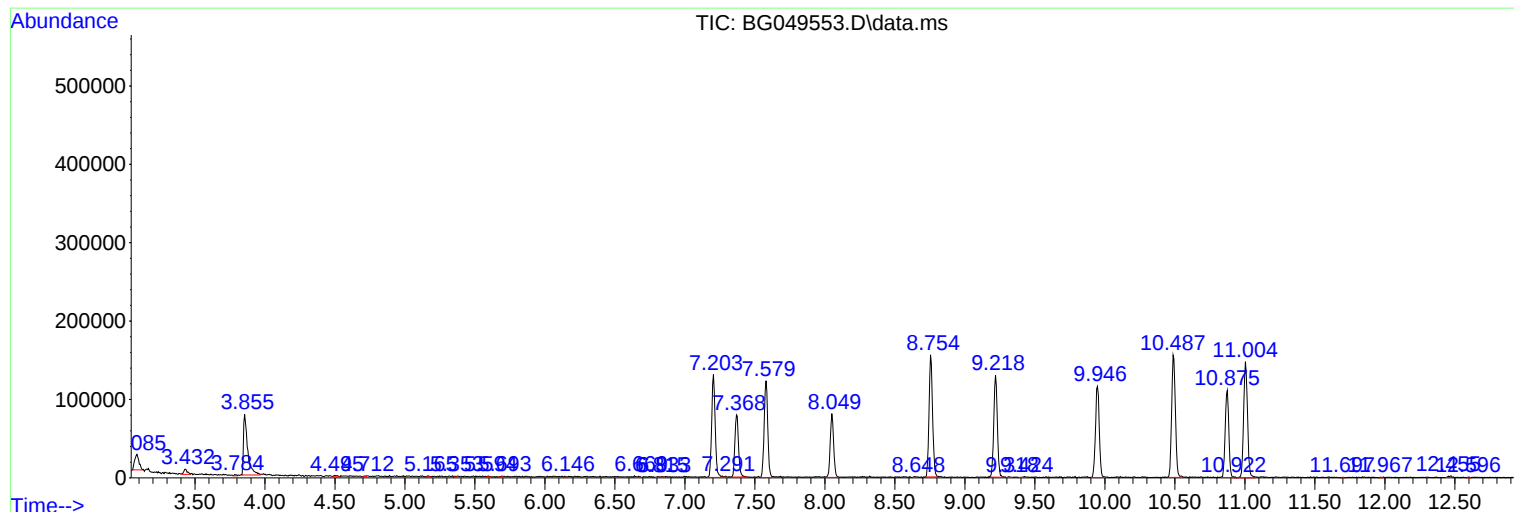
Sum of corrected areas: 8311797

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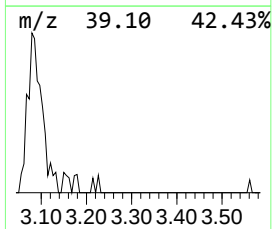
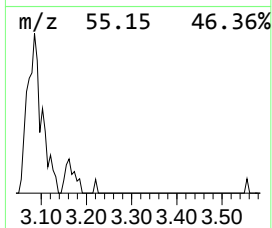
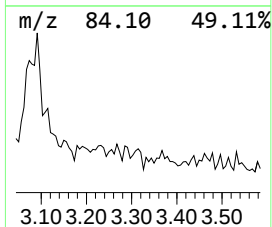
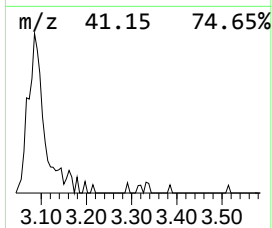
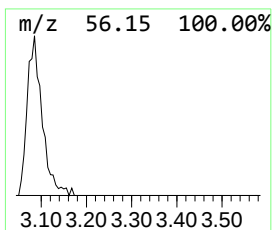
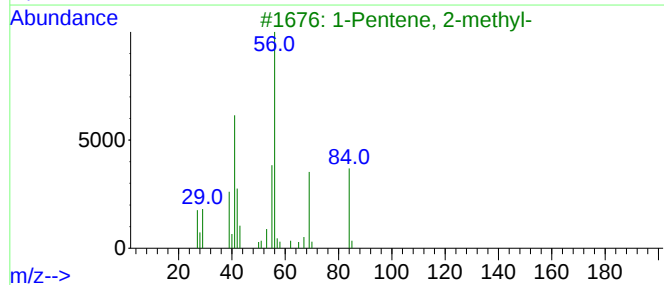
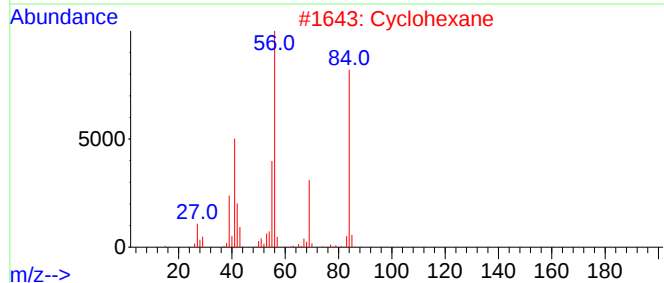
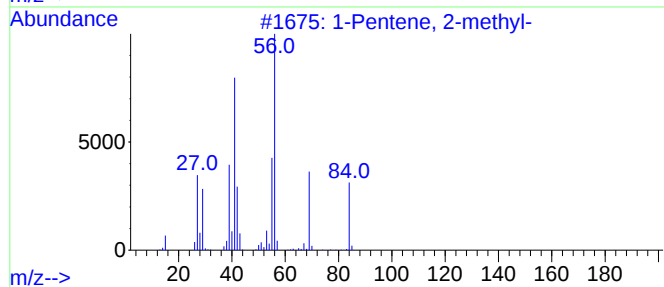
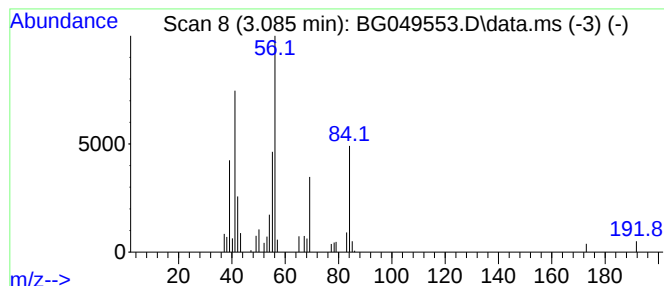
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG072421.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.085	6.06 ng/ul	41793	1,4-Dichlorobenzene-d4	8.049

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
2		Cyclohexane	84	C6H12	000110-82-7	58
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	58
4		2-Amino-oxazole	84	C3H4N2O	004570-45-0	53
5		Cyclohexane	84	C6H12	000110-82-7	53



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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: S...	3.085	6.1	ng/ul	41793	1	8.049	137926	20.0