

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010450.D
 Acq On : 20 May 2022 23:40
 Operator : CG/JU
 Sample : N2843-19
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4J9

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_P\Methods\SFAM-EPA-BP051322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010450.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	40	43	52	rVB	28767	47116	1.20%	0.118%
2	3.634	91	93	103	rVB3	5850	13196	0.34%	0.033%
3	3.764	112	115	134	rBV	173817	351249	8.92%	0.883%
4	3.987	149	153	160	rVB3	10851	19123	0.49%	0.048%
5	4.081	166	169	174	rVB2	8570	11701	0.30%	0.029%
6	4.999	319	325	333	rBV	38669	66235	1.68%	0.166%
7	7.058	666	675	688	rBV	524425	923232	23.44%	2.320%
8	7.222	695	703	715	rBV	443035	743341	18.88%	1.868%
9	7.311	715	718	726	rVB	7976	17461	0.44%	0.044%
10	7.422	730	737	751	rBV	559644	925259	23.50%	2.325%
11	7.516	751	753	757	rVB4	5589	7242	0.18%	0.018%
12	7.893	809	817	826	rBV	521074	868964	22.07%	2.184%
13	7.963	827	829	838	rVB8	7006	11731	0.30%	0.029%
14	8.599	928	937	952	rBV	652923	1123651	28.53%	2.824%
15	8.716	952	957	965	rVB	24821	46101	1.17%	0.116%
16	9.046	1006	1013	1031	rBV	577452	1031692	26.20%	2.593%
17	9.181	1031	1036	1039	rVV7	4608	8573	0.22%	0.022%
18	9.210	1039	1041	1050	rVB10	4079	8722	0.22%	0.022%
19	9.487	1084	1088	1098	rVB3	14105	24821	0.63%	0.062%
20	9.775	1129	1137	1145	rBV	479562	837218	21.26%	2.104%
21	10.134	1196	1198	1205	rVB8	3896	6726	0.17%	0.017%
22	10.310	1221	1228	1243	rBV	651575	1185285	30.10%	2.979%
23	10.693	1284	1293	1308	rBV	760990	1342431	34.09%	3.374%
24	10.828	1309	1316	1339	rVV	624192	1224134	31.08%	3.076%
25	11.734	1465	1470	1474	rBV8	2704	4510	0.11%	0.011%
26	12.128	1531	1537	1543	rBV2	8147	14702	0.37%	0.037%
27	12.199	1546	1549	1555	rBV8	2993	6009	0.15%	0.015%
28	12.281	1559	1563	1571	rVV4	12096	23777	0.60%	0.060%
29	12.928	1670	1673	1679	rVB7	5197	7609	0.19%	0.019%
30	13.016	1682	1688	1692	rBV9	3533	6263	0.16%	0.016%
31	13.075	1692	1698	1702	rVV9	5062	8614	0.22%	0.022%
32	13.134	1705	1708	1714	rVB6	5730	9962	0.25%	0.025%
33	13.357	1741	1746	1752	rVB	23859	32447	0.82%	0.082%
34	13.593	1783	1786	1791	rBV7	2291	4739	0.12%	0.012%
35	13.934	1836	1844	1855	rBV	1367446	2007532	50.98%	5.045%
36	14.057	1862	1865	1868	rVB5	5449	6293	0.16%	0.016%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
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37	14.216	1882	1892	1905	rBV	1510729	2277602	57.84%	5.724%
38	14.351	1911	1915	1921	rVB8	5787	9411	0.24%	0.024%
39	14.428	1924	1928	1934	rBV7	9758	16123	0.41%	0.041%
40	14.522	1937	1944	1957	rBV	1263307	1863687	47.32%	4.684%
41	14.722	1970	1978	1998	rBV	337737	760715	19.32%	1.912%
42	14.945	2012	2016	2024	rVB7	17749	31329	0.80%	0.079%
43	15.328	2077	2081	2086	rBV3	8552	16282	0.41%	0.041%
44	15.381	2086	2090	2093	rVV5	6144	8924	0.23%	0.022%
45	15.516	2106	2113	2128	rBV	2035035	2913328	73.98%	7.321%
46	15.634	2128	2133	2148	rVV	501127	780404	19.82%	1.961%
47	15.757	2150	2154	2162	rVB2	27290	45016	1.14%	0.113%
48	15.957	2186	2188	2192	rBV4	4722	6901	0.18%	0.017%
49	16.010	2192	2197	2202	rVB2	46699	64303	1.63%	0.162%
50	16.240	2235	2236	2241	rVV5	4587	5090	0.13%	0.013%
51	16.304	2243	2247	2251	rVB6	7697	11189	0.28%	0.028%
52	16.422	2265	2267	2272	rVB6	3630	4185	0.11%	0.011%
53	16.675	2306	2310	2314	rVB6	4146	5841	0.15%	0.015%
54	16.822	2332	2335	2337	rVB3	4480	4550	0.12%	0.011%
55	16.951	2355	2357	2361	rVB4	5256	5224	0.13%	0.013%
56	17.198	2392	2399	2403	rBV10	6466	12795	0.32%	0.032%
57	17.275	2403	2412	2422	rBV	1570841	2264182	57.49%	5.690%
58	17.375	2422	2429	2434	rVV	2179684	3187709	80.95%	8.011%
59	17.416	2434	2436	2442	rVV	104441	134405	3.41%	0.338%
60	17.475	2442	2446	2453	rVB7	23156	43257	1.10%	0.109%
61	17.945	2521	2526	2535	rVB	123496	158608	4.03%	0.399%
62	18.098	2549	2552	2556	rBV6	8727	11248	0.29%	0.028%
63	18.157	2558	2562	2570	rBV3	56575	85863	2.18%	0.216%
64	18.634	2640	2643	2648	rBV7	5689	10003	0.25%	0.025%
65	19.075	2714	2718	2726	rVB3	70478	82541	2.10%	0.207%
66	19.181	2732	2736	2744	rBV4	37229	57721	1.47%	0.145%
67	19.251	2744	2748	2751	rBV	28407	43157	1.10%	0.108%
68	19.392	2768	2772	2776	rBV6	26866	41613	1.06%	0.105%
69	19.533	2794	2796	2798	rBV3	10751	11217	0.28%	0.028%
70	19.604	2802	2808	2821	rBV	2888837	3938066	100.00%	9.897%
71	20.616	2977	2980	2983	rVB2	23167	21319	0.54%	0.054%
72	21.063	3052	3056	3063	rBV	208703	352372	8.95%	0.886%
73	21.357	3101	3106	3120	rBV	1703309	2309466	58.64%	5.804%
74	23.033	3388	3391	3397	rVB3	104035	161214	4.09%	0.405%
75	23.621	3484	3491	3497	rBV2	1516974	3029311	76.92%	7.613%
76	23.774	3511	3517	3533	rVB2	935557	2027751	51.49%	5.096%

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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_P\Methods\SFAM-EPA-BP051322.M
Title : SVOA CALIBRATION

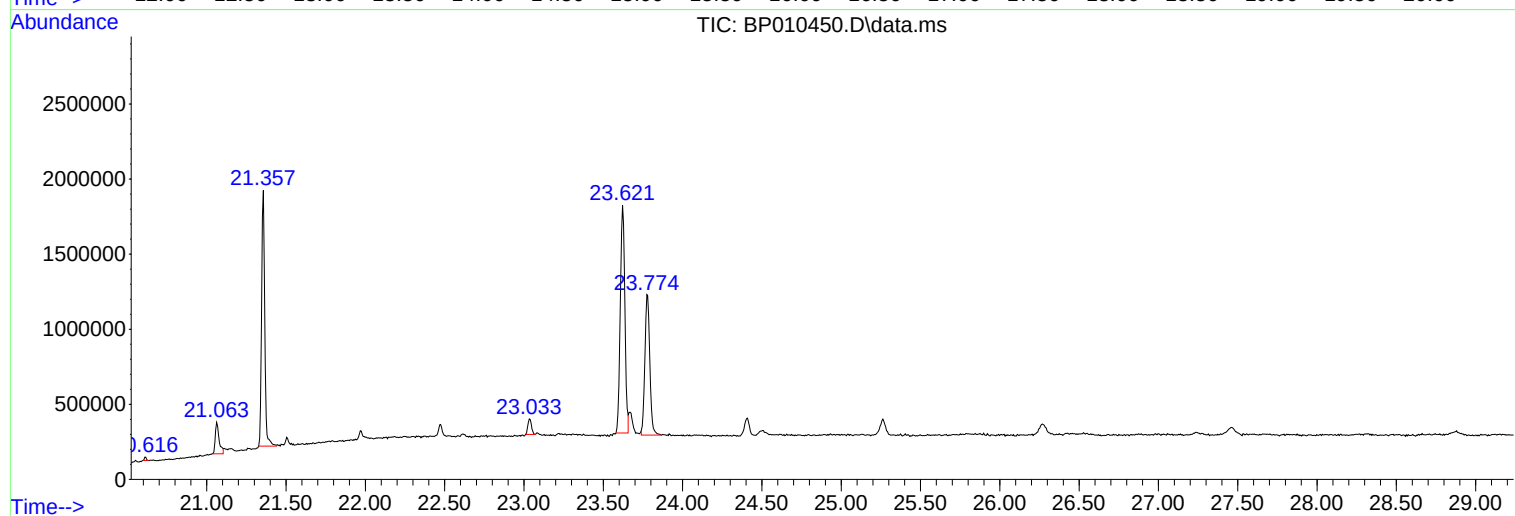
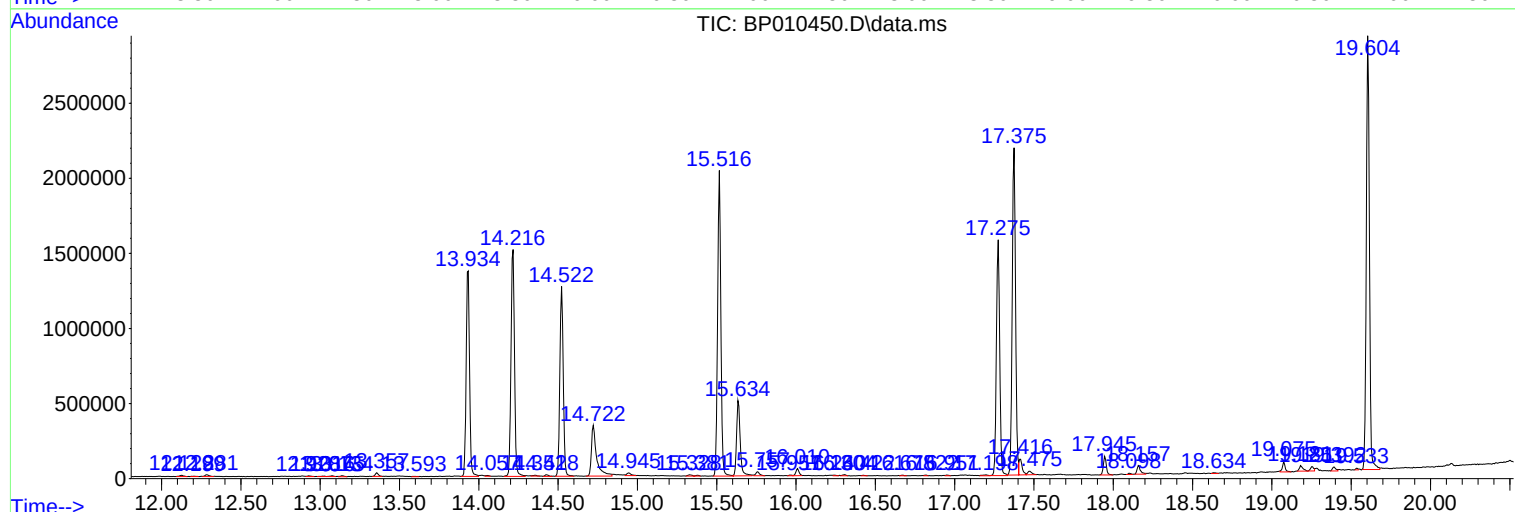
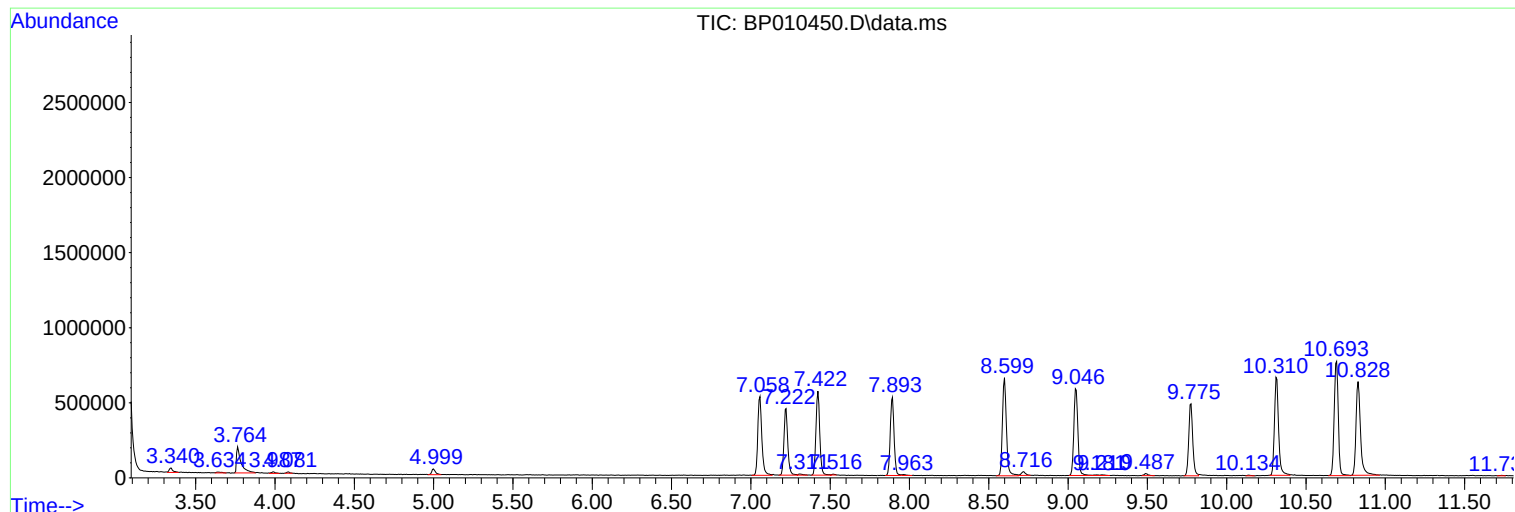
Sum of corrected areas: 39791583

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ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



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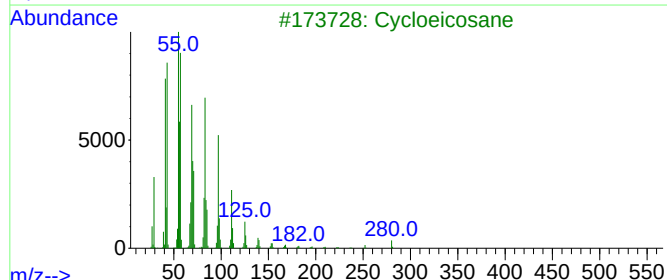
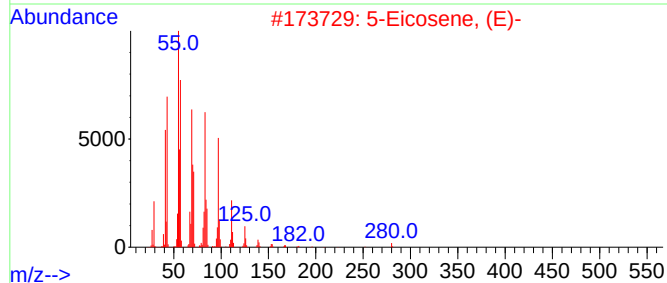
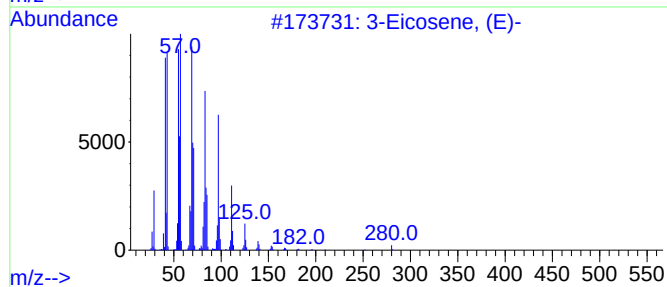
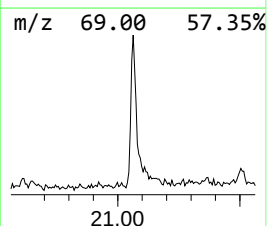
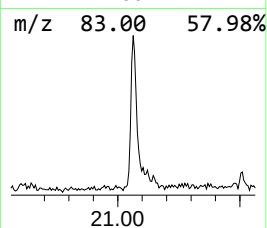
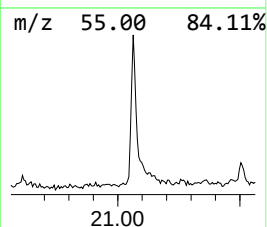
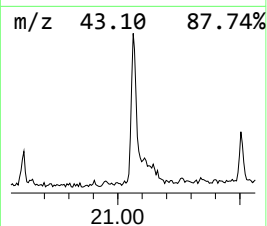
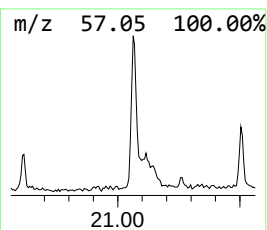
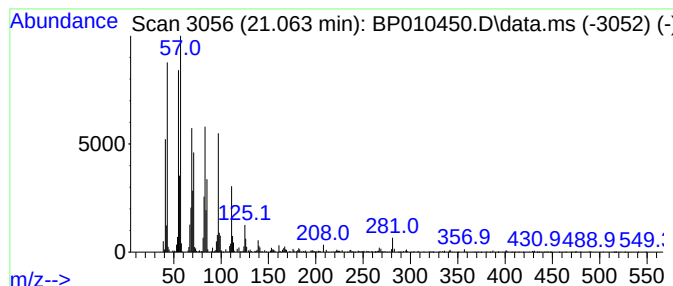
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.063	3.05 ng/ul	352372	Chrysene-d12	21.357

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-		280	C20H40	074685-33-9	92
2	5-Eicosene, (E)-		280	C20H40	074685-30-6	92
3	Cycloeicosane		280	C20H40	000296-56-0	91
4	1-Docosene		308	C22H44	001599-67-3	91
5	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	21.063	3.0	ng/u1	352372	5	21.357	2309470	20.0