

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120822\
 Data File : BP012993.D
 Acq On : 08 Dec 2022 19:57
 Operator : CG/JU
 Sample : N5894-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA8

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012993.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.357	25	29	37	rVB	136925	198909	0.96%	0.112%
2	3.787	99	102	118	rBV	589285	958744	4.62%	0.538%
3	4.016	136	141	148	rVB	75602	126819	0.61%	0.071%
4	4.116	154	158	164	rVB	49077	71877	0.35%	0.040%
5	4.493	220	222	227	rVB6	16469	23235	0.11%	0.013%
6	5.069	315	320	326	rBV3	14902	23366	0.11%	0.013%
7	6.046	480	486	491	rBV	16412	26728	0.13%	0.015%
8	7.128	664	670	681	rVB	638053	1013689	4.89%	0.569%
9	7.298	692	699	709	rBV	2384391	3675431	17.73%	2.062%
10	7.504	726	734	743	rBV	2214313	3490098	16.83%	1.958%
11	7.593	743	749	758	rVB10	16577	33522	0.16%	0.019%
12	7.975	807	814	820	rBV	2067482	3220893	15.53%	1.807%
13	8.034	820	824	835	rVB	425761	633538	3.06%	0.355%
14	8.681	927	934	945	rBV	1889366	2948011	14.22%	1.654%
15	8.804	950	955	966	rVB	31678	60841	0.29%	0.034%
16	9.075	995	1001	1005	rBV6	10678	23124	0.11%	0.013%
17	9.140	1005	1012	1022	rVV	2708928	4449306	21.46%	2.496%
18	9.245	1026	1030	1036	rVV	43069	71289	0.34%	0.040%
19	9.551	1078	1082	1088	rVB4	23791	41612	0.20%	0.023%
20	9.869	1128	1136	1149	rBV	2149402	3541905	17.08%	1.987%
21	10.004	1153	1159	1167	rVV	60000	121889	0.59%	0.068%
22	10.087	1170	1173	1179	rVB4	21842	33055	0.16%	0.019%
23	10.404	1219	1227	1239	rBV	3252930	5388892	25.99%	3.024%
24	10.563	1249	1254	1262	rBV	47231	85743	0.41%	0.048%
25	10.792	1283	1293	1301	rBV	2900665	4941493	23.83%	2.773%
26	10.922	1308	1315	1330	rBV	2230977	3849505	18.57%	2.160%
27	11.445	1400	1404	1411	rBV2	18154	30865	0.15%	0.017%
28	11.569	1420	1425	1432	rBV2	45599	86792	0.42%	0.049%
29	12.369	1556	1561	1568	rVB2	70728	115162	0.56%	0.065%
30	13.233	1701	1708	1713	rBV3	17768	35156	0.17%	0.020%
31	13.504	1748	1754	1762	rBV2	37814	71846	0.35%	0.040%
32	13.580	1763	1767	1774	rVB5	17473	35168	0.17%	0.020%
33	13.928	1820	1826	1835	rVB5	12973	29835	0.14%	0.017%
34	14.022	1835	1842	1857	rVV	6861007	9682386	46.70%	5.433%
35	14.133	1857	1861	1868	rVB10	13920	27149	0.13%	0.015%
36	14.257	1876	1882	1884	rBV4	44691	77350	0.37%	0.043%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	14.310	1884	1891	1903	rVB	7971453	11508652	55.51%	6.457%
38	14.498	1920	1923	1927	rVB3	18920	26278	0.13%	0.015%
39	14.616	1932	1943	1951	rBV	5001871	7260572	35.02%	4.074%
40	14.698	1955	1957	1965	rVB8	12396	23847	0.12%	0.013%
41	14.798	1965	1974	1993	rBV	510546	898020	4.33%	0.504%
42	15.022	2008	2012	2017	rVB7	39410	64781	0.31%	0.036%
43	15.245	2045	2050	2054	rBV	34624	52345	0.25%	0.029%
44	15.404	2072	2077	2084	rBV2	67423	130809	0.63%	0.073%
45	15.451	2084	2085	2089	rVB3	25437	22797	0.11%	0.013%
46	15.610	2104	2112	2124	rBV	10238297	15096545	72.81%	8.470%
47	15.722	2124	2131	2144	rVV	3457729	4651180	22.43%	2.610%
48	15.845	2147	2152	2158	rVB3	55478	93495	0.45%	0.052%
49	15.969	2168	2173	2176	rVV5	16498	22723	0.11%	0.013%
50	16.010	2176	2180	2185	rVB5	23904	32887	0.16%	0.018%
51	16.063	2185	2189	2194	rBV5	14967	27285	0.13%	0.015%
52	16.145	2200	2203	2206	rBV5	16078	24022	0.12%	0.013%
53	16.245	2216	2220	2223	rBV4	19336	25165	0.12%	0.014%
54	16.286	2223	2227	2229	rVV2	31956	42772	0.21%	0.024%
55	16.316	2229	2232	2240	rVV9	32624	68259	0.33%	0.038%
56	16.392	2242	2245	2250	rVB5	26303	33948	0.16%	0.019%
57	16.498	2259	2263	2267	rVB5	28132	36154	0.17%	0.020%
58	16.833	2313	2320	2327	rBV5	28291	58704	0.28%	0.033%
59	17.116	2362	2368	2371	rBV4	34395	53870	0.26%	0.030%
60	17.163	2372	2376	2384	rVB2	26428	54824	0.26%	0.031%
61	17.274	2391	2395	2400	rVB2	67786	87095	0.42%	0.049%
62	17.369	2404	2411	2418	rBV	6424847	8974062	43.28%	5.035%
63	17.469	2422	2428	2439	rVV	11767690	17004719	82.02%	9.541%
64	17.551	2440	2442	2448	rVB5	46083	69404	0.33%	0.039%
65	17.651	2455	2459	2468	rBV	82812	143711	0.69%	0.081%
66	17.745	2471	2475	2478	rBV5	28984	36006	0.17%	0.020%
67	18.027	2519	2523	2527	rVB3	107882	136396	0.66%	0.077%
68	18.233	2554	2558	2564	rBV	205177	272282	1.31%	0.153%
69	18.310	2568	2571	2581	rVB	149639	200544	0.97%	0.113%
70	18.651	2626	2629	2634	rBV5	39200	51951	0.25%	0.029%
71	18.815	2655	2657	2661	rBV4	26950	27266	0.13%	0.015%
72	18.980	2682	2685	2689	rVB	43846	51696	0.25%	0.029%
73	19.063	2696	2699	2704	rBV3	101125	115215	0.56%	0.065%
74	19.274	2731	2735	2739	rBV2	167927	225985	1.09%	0.127%
75	19.339	2742	2746	2750	rBV	156233	227621	1.10%	0.128%
76	19.468	2764	2768	2777	rBV3	87864	158390	0.76%	0.089%

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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_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

77	19.698	2801	2807	2823	rBV	15460599	20733142	100.00%	11.633%
78	21.145	3049	3053	3058	rBV2	265314	402587	1.94%	0.226%
79	21.451	3100	3105	3115	rBV	7212869	9635369	46.47%	5.406%
80	23.162	3392	3396	3405	rVB	329028	539382	2.60%	0.303%
81	23.786	3494	3502	3515	rVB2	9325339	20245880	97.65%	11.359%
82	23.945	3522	3529	3541	rVB	4644175	9334886	45.02%	5.238%

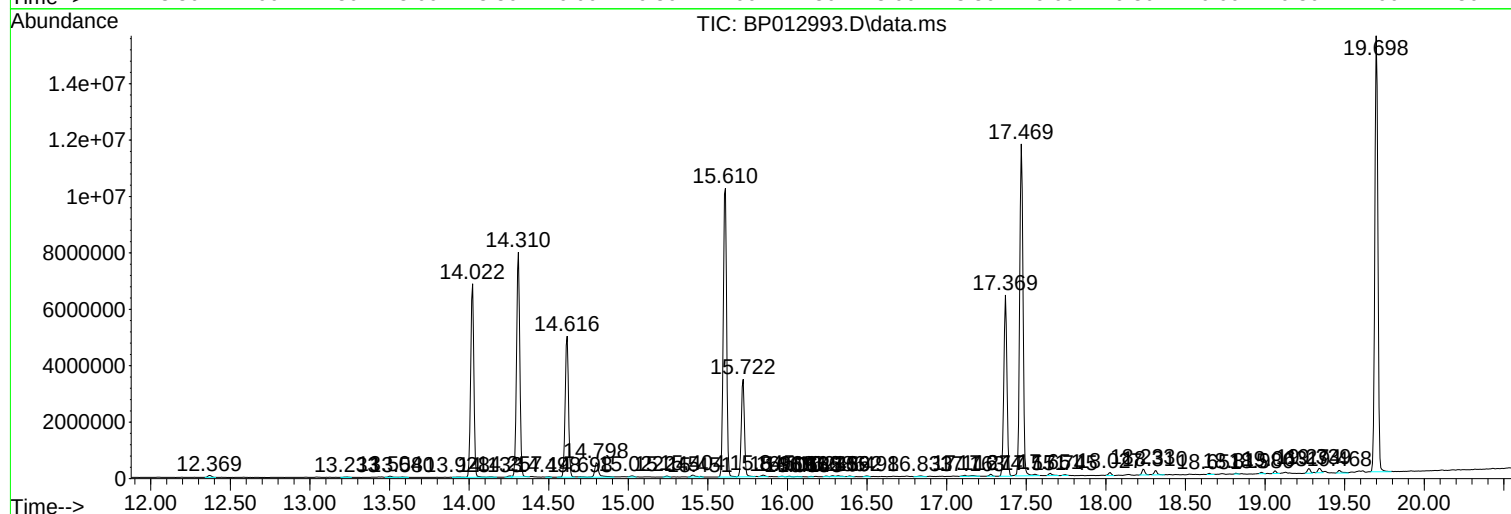
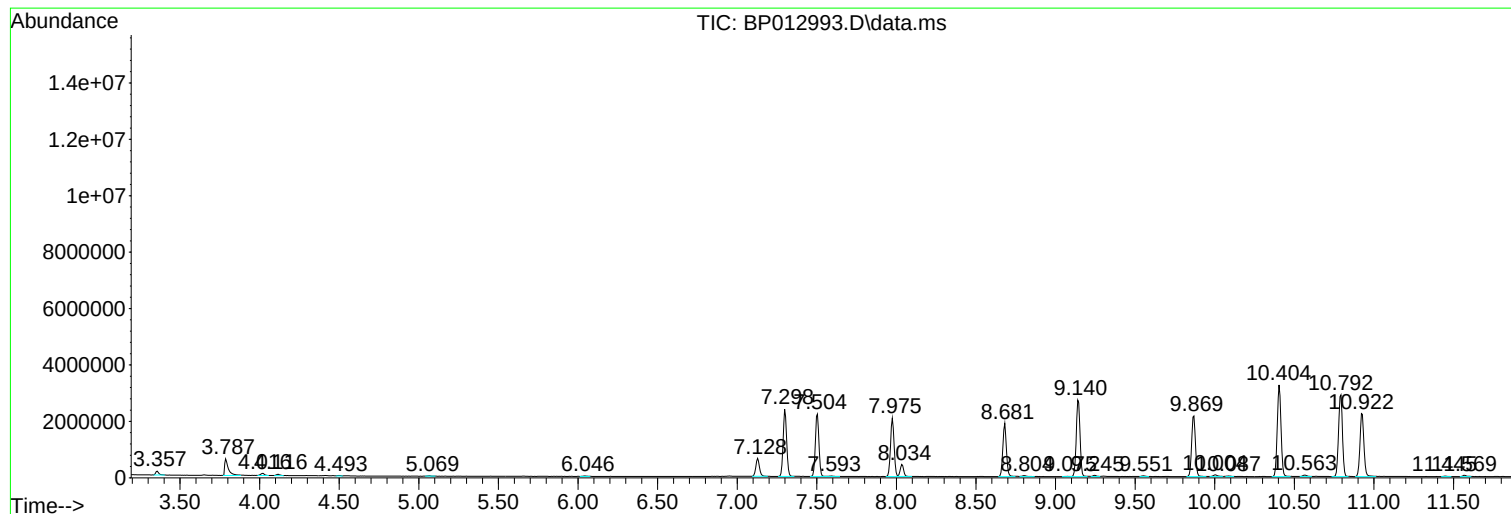
Sum of corrected areas: 178228746

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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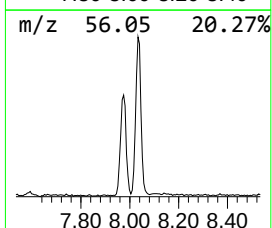
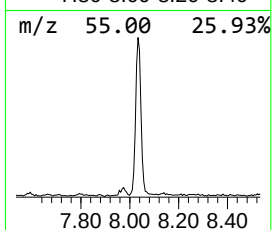
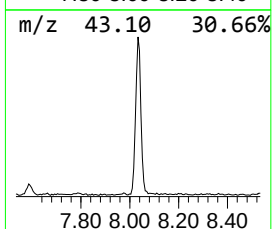
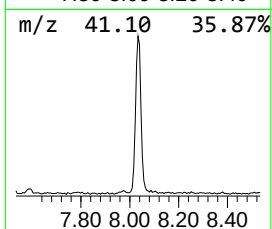
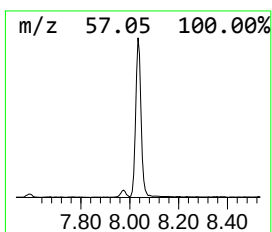
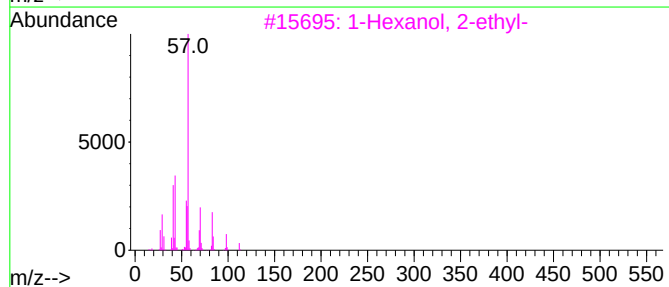
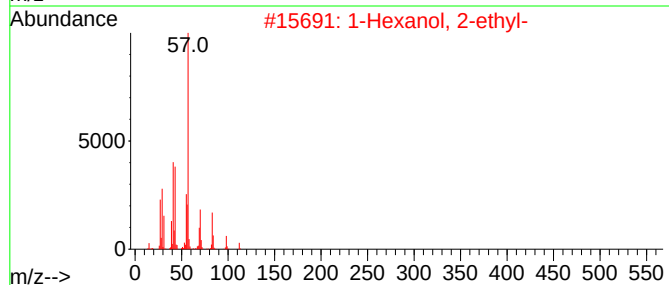
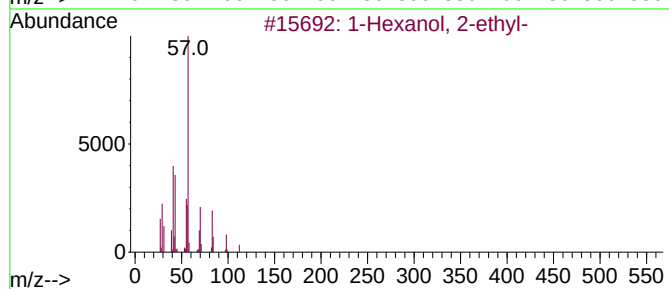
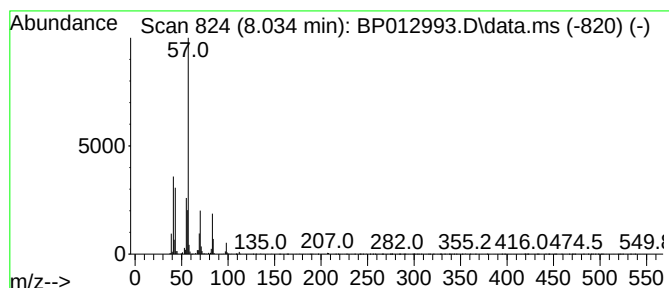
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.034	3.93 ng/ul	633538	1,4-Dichlorobenzene-d4	7.975

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	8.034	3.9	ng/u1	633538	1	7.975	3220890	20.0