

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
 Data File : BP021434.D
 Acq On : 08 Aug 2024 01:45
 Operator : CG/JU
 Sample : P3401-10
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1GF5

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021434.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.152	24	28	33	rBV	32878	41791	0.85%	0.107%
2	3.605	102	105	124	rBV	31921	94815	1.93%	0.244%
3	3.864	146	149	155	rBV3	7839	13356	0.27%	0.034%
4	5.717	459	464	469	rBV8	5087	9420	0.19%	0.024%
5	6.875	651	661	671	rBV2	65596	231184	4.70%	0.594%
6	6.964	671	676	703	rVV	390839	865952	17.59%	2.227%
7	7.169	704	711	739	rVB	406205	974083	19.78%	2.505%
8	7.605	778	785	794	rBV	458485	881165	17.90%	2.266%
9	7.675	794	797	813	rVB3	46956	139468	2.83%	0.359%
10	7.881	830	832	837	rVB5	7031	10418	0.21%	0.027%
11	8.363	907	914	926	rBV	196240	601246	12.21%	1.546%
12	8.463	927	931	945	rVB	47849	118873	2.41%	0.306%
13	8.705	966	972	978	rBV2	21462	46852	0.95%	0.120%
14	8.781	978	985	1004	rVV	408492	822934	16.71%	2.116%
15	9.493	1097	1106	1127	rBV	277991	717799	14.58%	1.846%
16	9.775	1149	1154	1160	rVB10	2600	4997	0.10%	0.013%
17	10.058	1195	1202	1219	rBV	270202	960034	19.50%	2.468%
18	10.316	1240	1246	1249	rBV6	13077	23458	0.48%	0.060%
19	10.375	1249	1256	1264	rBV	560348	1032889	20.98%	2.656%
20	10.563	1280	1288	1315	rBV	208443	728306	14.79%	1.873%
21	11.128	1381	1384	1397	rVB8	5337	15417	0.31%	0.040%
22	12.010	1528	1534	1544	rBV3	7539	19594	0.40%	0.050%
23	12.834	1669	1674	1683	rVB5	10849	20221	0.41%	0.052%
24	13.110	1716	1721	1730	rBV4	7087	18108	0.37%	0.047%
25	13.475	1779	1783	1791	rBV10	3585	6187	0.13%	0.016%
26	13.669	1809	1816	1828	rBV	1064050	1790792	36.37%	4.604%
27	13.940	1853	1862	1888	rBV2	1176065	2138036	43.42%	5.497%
28	14.122	1888	1893	1899	rVB10	3271	7904	0.16%	0.020%
29	14.246	1904	1914	1925	rBV	976057	1580068	32.09%	4.063%
30	14.328	1925	1928	1940	rVB3	16330	30150	0.61%	0.078%
31	14.704	1986	1992	1993	rBV4	15686	20438	0.42%	0.053%
32	14.934	2027	2031	2033	rBV5	5178	7689	0.16%	0.020%
33	15.045	2046	2050	2054	rVB	8723	11922	0.24%	0.031%
34	15.110	2058	2061	2067	rVB3	4962	7928	0.16%	0.020%
35	15.269	2081	2088	2106	rBV	1547732	2762890	56.11%	7.104%
36	15.481	2116	2124	2138	rBV2	228401	689913	14.01%	1.774%

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37	15.663	2151	2155	2164	rVB3	16362	34066	0.69%	0.088%
38	15.845	2182	2186	2191	rBV8	2891	5381	0.11%	0.014%
39	16.645	2319	2322	2329	rBV8	4900	10796	0.22%	0.028%
40	16.863	2357	2359	2365	rVB5	4654	6713	0.14%	0.017%
41	17.022	2381	2386	2388	rBV5	3614	6881	0.14%	0.018%
42	17.069	2388	2394	2405	rVV	1301970	1984695	40.31%	5.103%
43	17.181	2405	2413	2435	rVB	1738062	3202817	65.05%	8.235%
44	17.375	2442	2446	2451	rVB	11796	17608	0.36%	0.045%
45	17.757	2504	2511	2519	rBV2	252028	354510	7.20%	0.912%
46	18.010	2549	2554	2562	rBV3	33990	70944	1.44%	0.182%
47	18.081	2563	2566	2571	rVB	14745	25803	0.52%	0.066%
48	18.216	2587	2589	2593	rBV5	4883	6177	0.13%	0.016%
49	18.928	2707	2710	2711	rBV3	11456	11657	0.24%	0.030%
50	19.104	2736	2740	2745	rBV2	27689	45128	0.92%	0.116%
51	19.198	2752	2756	2759	rBV2	30204	42155	0.86%	0.108%
52	19.345	2777	2781	2786	rBV7	31052	53835	1.09%	0.138%
53	19.569	2812	2819	2830	rBV	2812208	4773554	96.95%	12.274%
54	21.522	3145	3151	3165	rBV2	1439662	2787621	56.62%	7.168%
55	24.574	3661	3670	3688	rVB	1651084	4923744	100.00%	12.660%
56	24.804	3700	3709	3725	rVB	1104955	3081854	62.59%	7.924%

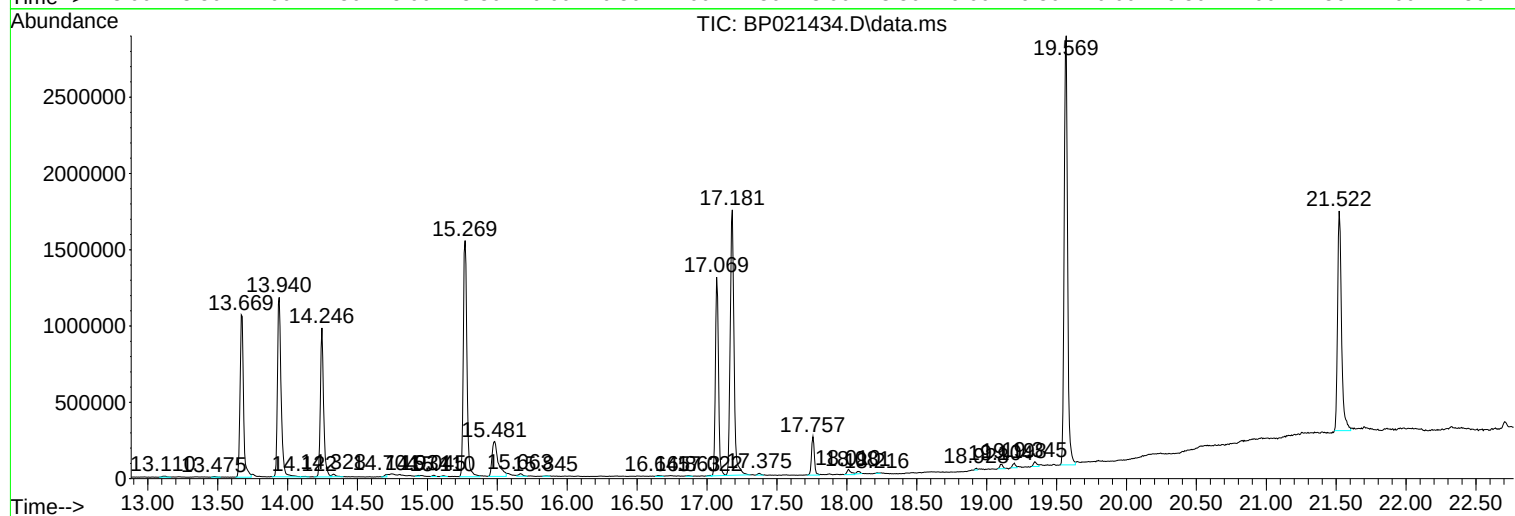
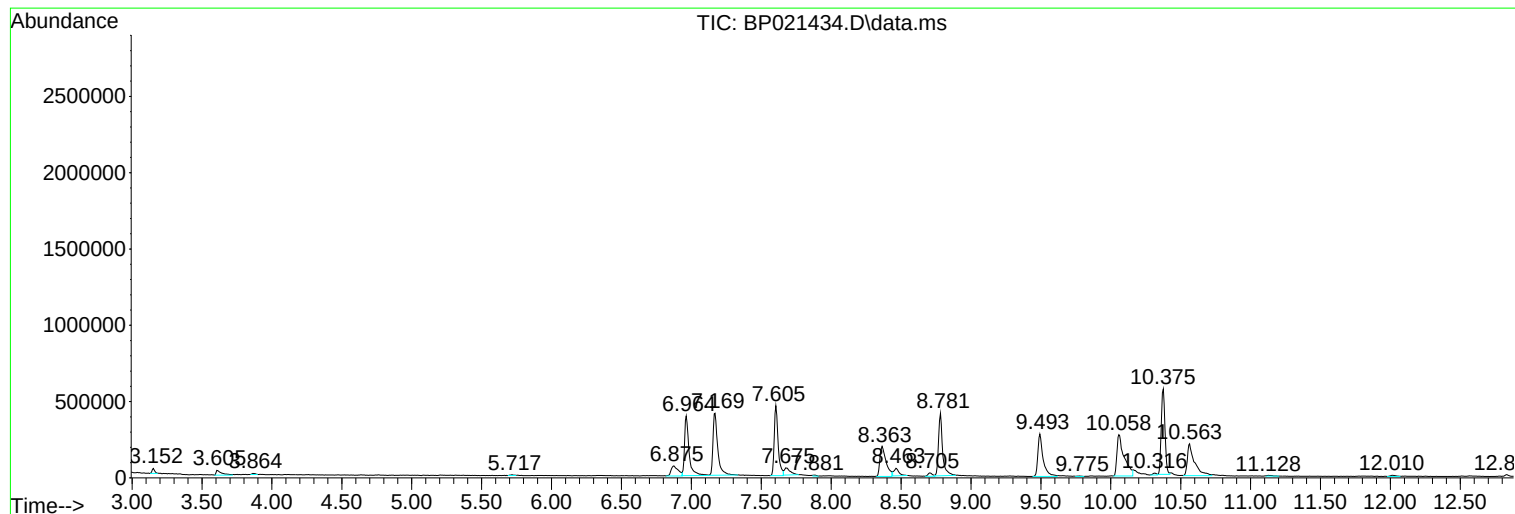
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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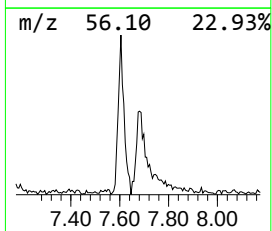
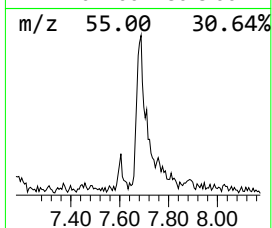
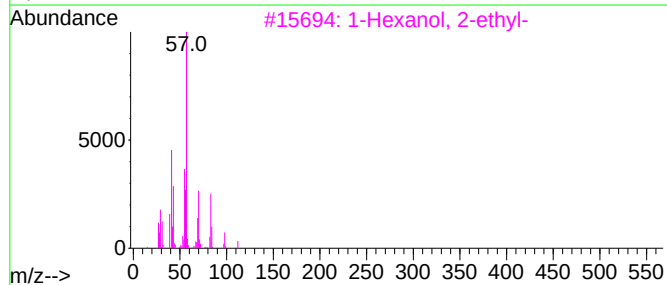
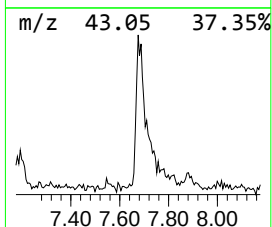
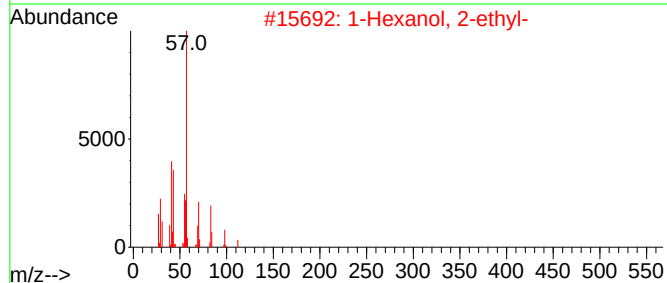
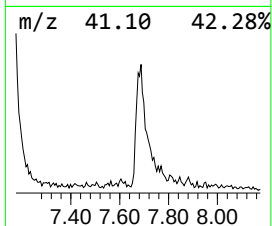
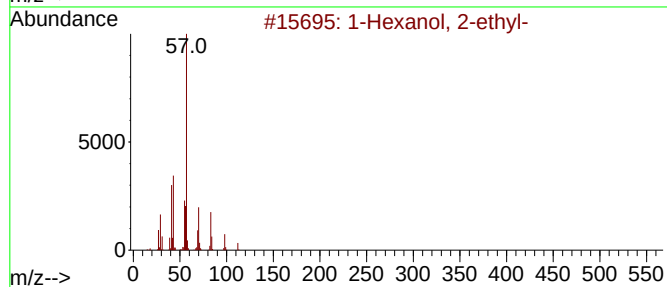
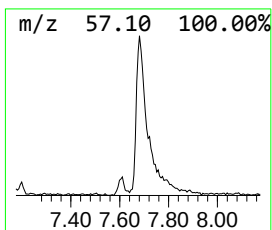
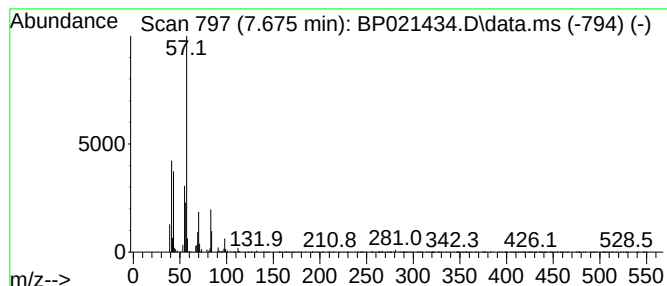
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.675	3.17 ng/ul	139468	1,4-Dichlorobenzene-d4	7.605

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	1-Decene, 2,4-dimethyl-			168	C12H24	055170-80-4	50
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	50



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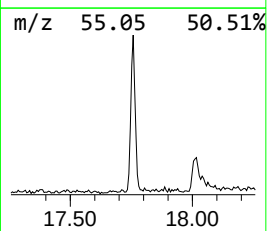
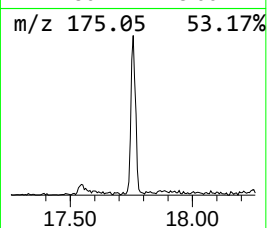
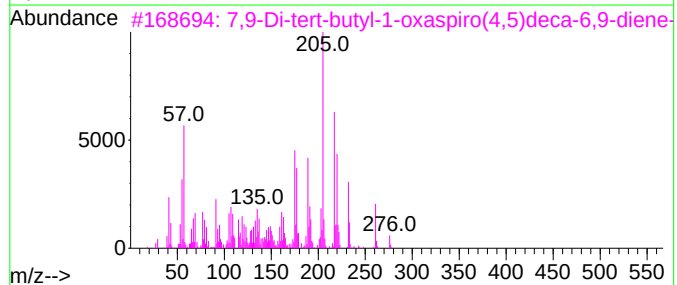
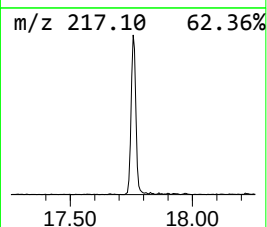
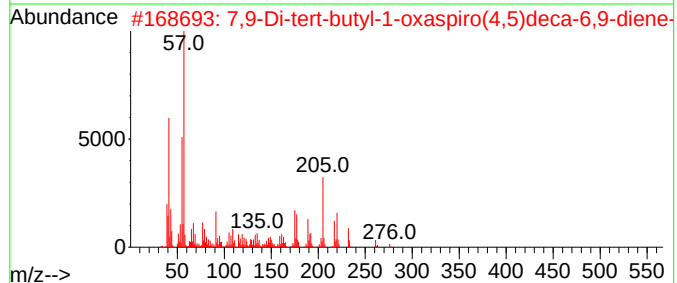
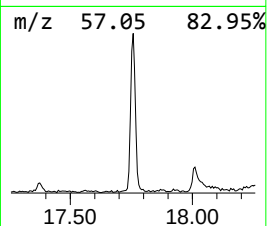
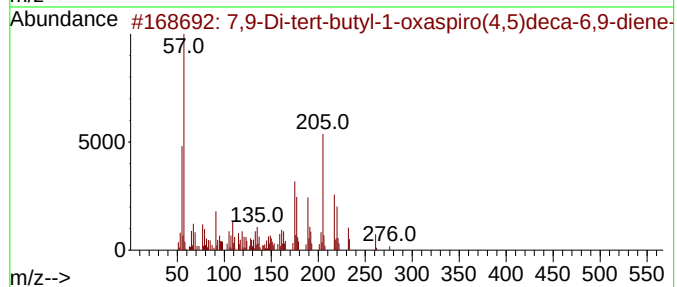
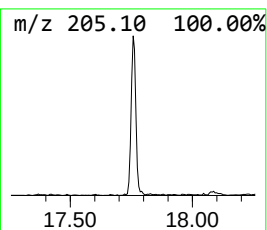
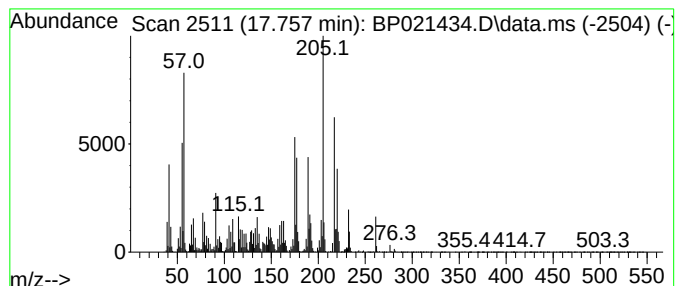
Instrument :
BNA_P
ClientSampleId :
E1GF5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.757	3.57 ng/ul	354510	Phenanthrene-d10	17.069	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	97
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	91
4	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	83
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	45



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	7.675	3.2	ng/ul	139468	1	7.605	881165	20.0
7,9-Di-tert-but...	17.757	3.6	ng/ul	354510	4	17.069	1984700	20.0