

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018167.D
 Acq On : 14 Nov 2023 21:47
 Operator : MA/JU
 Sample : 05337-13
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDJ8

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018167.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	9	rVB	60874	60354	2.41%	0.284%
2	3.164	9	13	17	rVB	15573	21311	0.85%	0.100%
3	3.622	87	91	97	rBV	21828	43084	1.72%	0.203%
4	3.775	115	117	124	rVB4	5584	8034	0.32%	0.038%
5	3.899	135	138	139	rBV3	5567	5333	0.21%	0.025%
6	4.417	222	226	231	rBV6	8064	9949	0.40%	0.047%
7	4.587	253	255	264	rBV8	2230	4407	0.18%	0.021%
8	5.481	402	407	409	rVB6	1733	2910	0.12%	0.014%
9	6.969	655	660	673	rVV	32983	81554	3.25%	0.384%
10	7.140	683	689	701	rBV	229295	375855	14.99%	1.771%
11	7.310	712	718	730	rBV	213856	383341	15.29%	1.806%
12	7.787	792	799	812	rBV	289545	492280	19.63%	2.320%
13	8.516	918	923	939	rBV	129964	264985	10.57%	1.249%
14	8.993	998	1004	1019	rBV	283573	530646	21.16%	2.501%
15	9.704	1118	1125	1139	rBV	235092	452295	18.04%	2.131%
16	10.234	1208	1215	1236	rBV	265772	579667	23.11%	2.732%
17	10.404	1243	1244	1249	rVV5	1797	2618	0.10%	0.012%
18	10.604	1270	1278	1290	rBV	398727	671914	26.79%	3.166%
19	10.787	1302	1309	1334	rBV	268364	632824	25.23%	2.982%
20	11.657	1449	1457	1464	rVB5	6678	13439	0.54%	0.063%
21	11.728	1467	1469	1475	rVB7	1707	2579	0.10%	0.012%
22	12.240	1551	1556	1557	rBV5	2990	4147	0.17%	0.020%
23	12.257	1557	1559	1563	rVB5	2422	2808	0.11%	0.013%
24	12.769	1640	1646	1649	rBV7	3322	4571	0.18%	0.022%
25	13.034	1684	1691	1692	rBV7	3278	3690	0.15%	0.017%
26	13.304	1731	1737	1743	rBV5	8796	14278	0.57%	0.067%
27	13.498	1765	1770	1773	rVB7	1935	2780	0.11%	0.013%
28	13.675	1795	1800	1805	rVB8	3432	6363	0.25%	0.030%
29	13.781	1812	1818	1819	rBV2	14503	20397	0.81%	0.096%
30	13.887	1831	1836	1857	rVB	845806	1191016	47.49%	5.612%
31	14.163	1877	1883	1893	rBV2	756189	1106657	44.13%	5.215%
32	14.463	1926	1934	1948	rBV	597755	893987	35.65%	4.213%
33	14.563	1948	1951	1955	rBV6	2334	3159	0.13%	0.015%
34	14.792	1985	1990	1994	rBV7	7895	15986	0.64%	0.075%
35	15.204	2055	2060	2066	rBV3	8601	15277	0.61%	0.072%
36	15.251	2066	2068	2071	rVB4	2575	3129	0.12%	0.015%

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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-BP111023.MA.M
 Title : SVOA CALIBRATION

37	15.304	2074	2077	2081	rVB6	3461	4756	0.19%	0.022%
38	15.469	2099	2105	2121	rBV	1123806	1579194	62.97%	7.442%
39	15.628	2126	2132	2145	rBV	258905	467495	18.64%	2.203%
40	15.716	2145	2147	2158	rVB8	13107	23189	0.92%	0.109%
41	15.886	2172	2176	2179	rBV	5965	6707	0.27%	0.032%
42	15.922	2181	2182	2186	rVV4	4212	4729	0.19%	0.022%
43	16.186	2222	2227	2233	rBV2	15115	21759	0.87%	0.103%
44	16.281	2237	2243	2247	rBV	41484	58900	2.35%	0.278%
45	16.345	2248	2254	2259	rVV3	61116	114697	4.57%	0.540%
46	16.404	2259	2264	2268	rVV2	37618	62295	2.48%	0.294%
47	16.445	2268	2271	2275	rVB2	23222	27412	1.09%	0.129%
48	16.504	2276	2281	2282	rBV3	10006	13426	0.54%	0.063%
49	16.545	2286	2288	2293	rVB3	10979	15238	0.61%	0.072%
50	16.604	2293	2298	2304	rBV4	29333	52726	2.10%	0.248%
51	16.675	2304	2310	2314	rBV	37823	52941	2.11%	0.249%
52	16.745	2319	2322	2328	rVB2	19055	28031	1.12%	0.132%
53	16.881	2340	2345	2347	rBV6	2373	3212	0.13%	0.015%
54	16.910	2347	2350	2357	rVV9	3588	4772	0.19%	0.022%
55	16.986	2360	2363	2365	rVV4	2521	3216	0.13%	0.015%
56	17.045	2369	2373	2377	rBV6	2948	4097	0.16%	0.019%
57	17.133	2386	2388	2391	rVB4	2715	2674	0.11%	0.013%
58	17.245	2400	2407	2418	rBV	820345	1135075	45.26%	5.349%
59	17.345	2418	2424	2441	rVV	1293536	1914532	76.34%	9.022%
60	17.504	2447	2451	2453	rBV5	4059	5691	0.23%	0.027%
61	17.586	2462	2465	2468	rBV5	4484	5429	0.22%	0.026%
62	17.675	2477	2480	2483	rBV5	4435	5128	0.20%	0.024%
63	17.886	2508	2516	2518	rBV9	10911	17392	0.69%	0.082%
64	17.910	2518	2520	2524	rVB4	3030	3063	0.12%	0.014%
65	18.069	2540	2547	2549	rBV7	2309	4937	0.20%	0.023%
66	18.098	2549	2552	2563	rVV7	15250	30566	1.22%	0.144%
67	18.192	2565	2568	2575	rVB	10045	13289	0.53%	0.063%
68	18.310	2584	2588	2591	rBV6	3373	5724	0.23%	0.027%
69	18.563	2626	2631	2635	rBV8	2357	4880	0.19%	0.023%
70	19.175	2731	2735	2740	rBV3	16624	26896	1.07%	0.127%
71	19.239	2741	2746	2753	rVV4	28993	61940	2.47%	0.292%
72	19.351	2762	2765	2775	rVB4	7811	14036	0.56%	0.066%
73	19.480	2783	2787	2790	rBV4	14863	16543	0.66%	0.078%
74	19.604	2802	2808	2817	rBV	1839986	2360334	94.12%	11.123%
75	20.510	2959	2962	2964	rBV4	11458	14467	0.58%	0.068%
76	21.380	3105	3110	3120	rBV	917735	1266146	50.49%	5.967%

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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-BP111023.MA.M
Title : SVOA CALIBRATION

77	23.686	3494	3502	3511	rBV	1291308	2507789	100.00%	11.818%
78	23.845	3523	3529	3543	rVB	602619	1327886	52.95%	6.257%

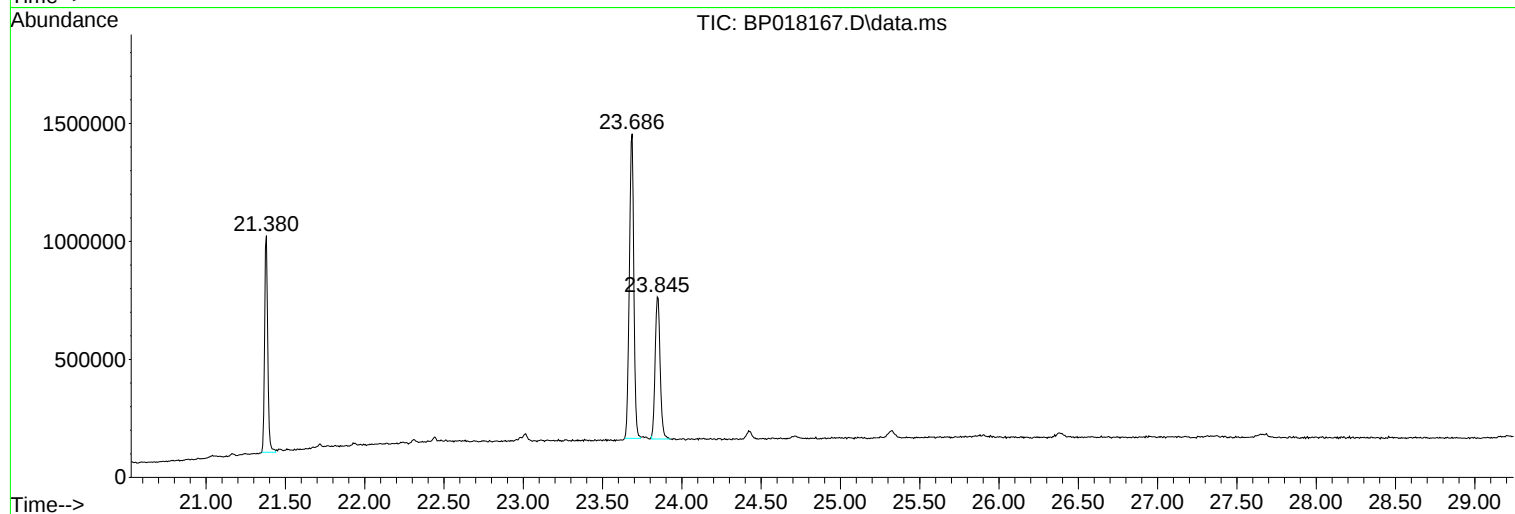
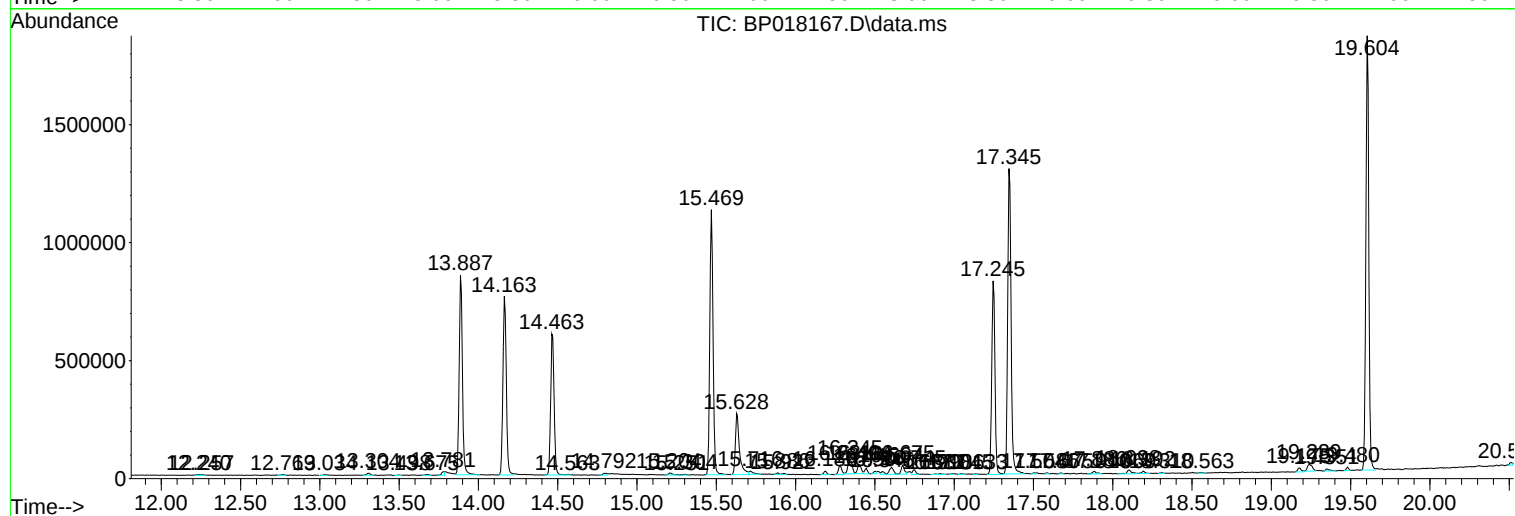
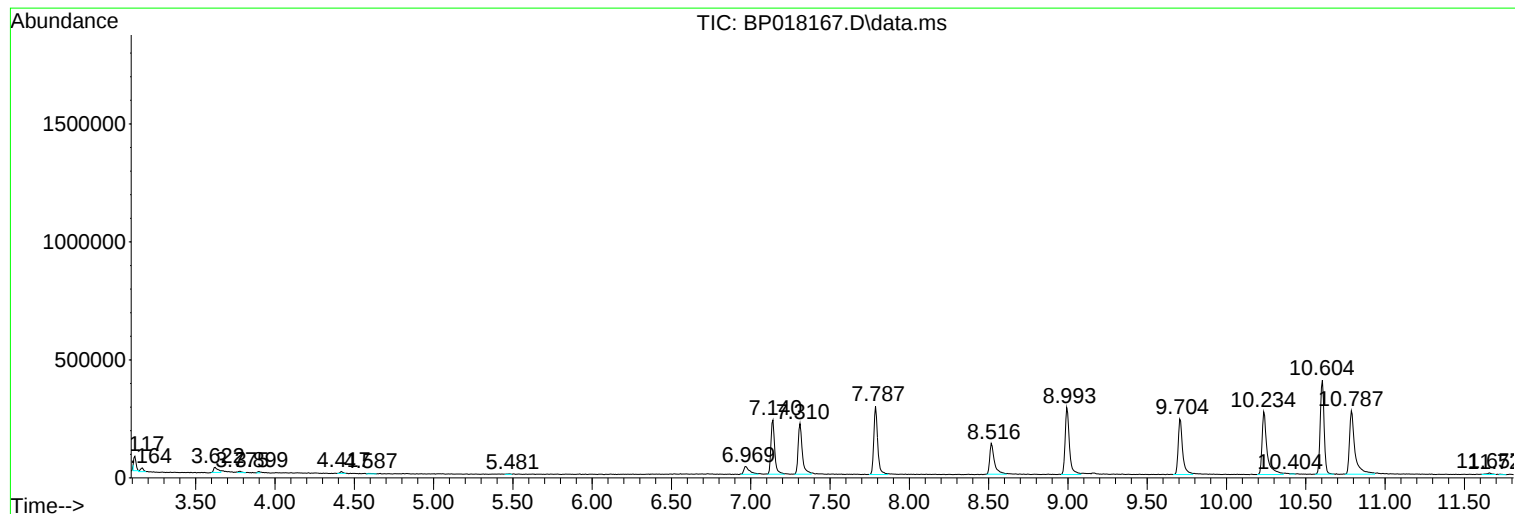
Sum of corrected areas: 21220833

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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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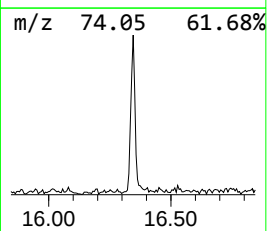
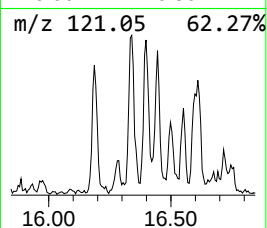
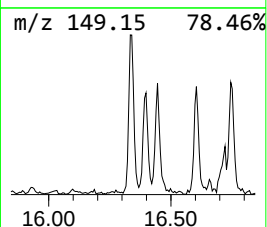
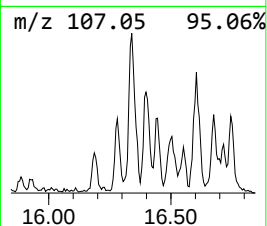
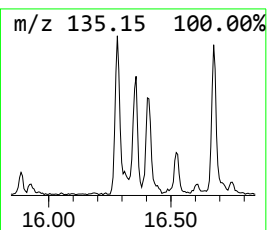
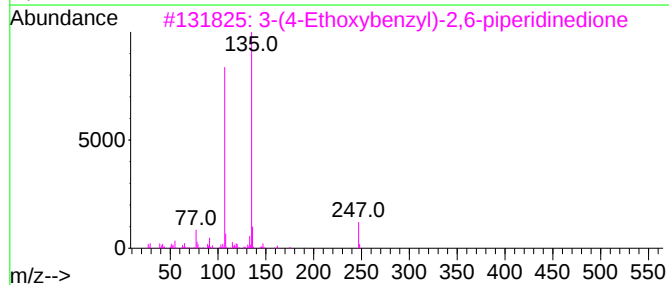
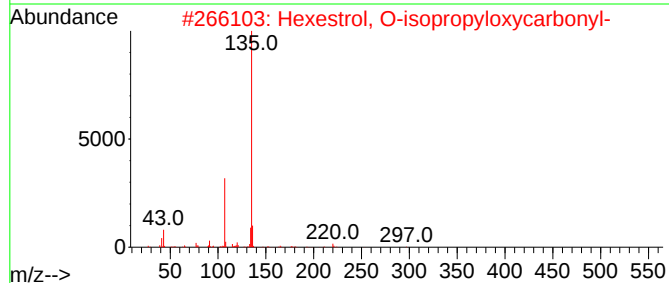
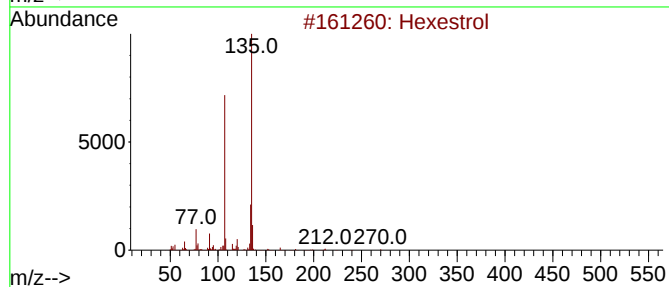
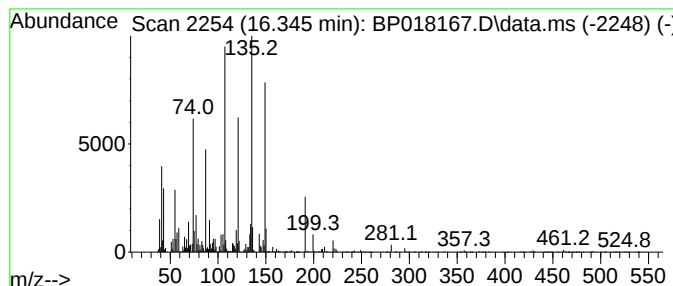
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	2.02 ng/ul	114697	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexestrol	270	C18H22O2	000084-16-2	41
2			Hexestrol, O-isopropylloxycarbonyl-	356	C22H28O4	1000487-68-4	35
3			3-(4-Ethoxybenzyl)-2,6-piperidin...	247	C14H17N03	051180-37-1	35
4			Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	30
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	22



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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
unknown-01	16.345	2.0	ng/u1	114697	4	17.245	1135080	20.0