

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041724\
 Data File : BN030981.D
 Acq On : 17 Apr 2024 17:27
 Operator : MA/JU
 Sample : P2060-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E1CY9

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_N\Methods\SFAM-EPA-BN032824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN030981.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	23717	29168	0.36%	0.050%
2	3.564	96	98	115	rBV	100202	182408	2.24%	0.316%
3	3.875	147	151	162	rVB2	11401	20850	0.26%	0.036%
4	5.211	373	378	382	rVB4	5504	9243	0.11%	0.016%
5	5.511	425	429	434	rVB4	5814	8984	0.11%	0.016%
6	5.752	466	470	475	rVB6	4400	8197	0.10%	0.014%
7	6.816	645	651	661	rVB	160931	259463	3.18%	0.449%
8	6.981	673	679	692	rBV	641019	1034715	12.68%	1.790%
9	7.175	704	712	723	rBV	590639	940727	11.53%	1.627%
10	7.640	785	791	797	rBV	599962	931644	11.42%	1.612%
11	7.705	797	802	811	rVB	150621	227378	2.79%	0.393%
12	8.346	904	911	921	rBV	521806	811806	9.95%	1.404%
13	8.463	927	931	937	rVB2	13793	21387	0.26%	0.037%
14	8.728	972	976	981	rBV5	6346	10671	0.13%	0.018%
15	8.799	981	988	999	rVV	774923	1272259	15.59%	2.201%
16	9.510	1102	1109	1125	rBV	524284	857699	10.51%	1.484%
17	10.046	1193	1200	1211	rBV	879440	1469414	18.01%	2.542%
18	10.210	1223	1228	1232	rVB6	9571	15557	0.19%	0.027%
19	10.422	1257	1264	1270	rBV	899577	1517450	18.60%	2.625%
20	10.463	1270	1271	1278	rVB3	34904	46234	0.57%	0.080%
21	10.563	1280	1288	1303	rBV	859334	1452404	17.80%	2.513%
22	11.081	1369	1376	1382	rVB4	5438	11319	0.14%	0.020%
23	12.016	1531	1535	1542	rBV	16339	24819	0.30%	0.043%
24	12.887	1678	1683	1689	rBV3	8517	12463	0.15%	0.022%
25	13.128	1718	1724	1730	rBV2	19783	28744	0.35%	0.050%
26	13.181	1730	1733	1739	rVB4	12470	17557	0.22%	0.030%
27	13.704	1815	1822	1835	rBV	2212622	3007331	36.85%	5.203%
28	13.975	1861	1868	1879	rBV	2300629	3419024	41.90%	5.915%
29	14.287	1914	1921	1932	rBV	1599290	2295101	28.13%	3.970%
30	14.498	1951	1957	1975	rBV	137363	265716	3.26%	0.460%
31	14.716	1988	1994	1998	rVB6	4391	8731	0.11%	0.015%
32	14.969	2031	2037	2042	rBV	20658	33920	0.42%	0.059%
33	15.098	2054	2059	2067	rVB4	16991	34795	0.43%	0.060%
34	15.281	2084	2090	2103	rBV	3083468	4493813	55.07%	7.774%
35	15.404	2105	2111	2121	rBV	821893	1077095	13.20%	1.863%
36	15.522	2127	2131	2136	rVB2	14444	22641	0.28%	0.039%

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 Title : SVOA CALIBRATION

37	15.575	2136	2140	2144	rVV5	7094	9503	0.12%	0.016%
38	15.651	2148	2153	2158	rBV3	6619	8915	0.11%	0.015%
39	15.898	2191	2195	2200	rVB7	6895	10013	0.12%	0.017%
40	16.069	2220	2224	2229	rBV4	9861	14045	0.17%	0.024%
41	17.033	2382	2388	2398	rVB	2122509	2944831	36.09%	5.094%
42	17.133	2398	2405	2420	rBV2	3704615	5303570	65.00%	9.175%
43	17.328	2434	2438	2443	rBV	17954	21611	0.26%	0.037%
44	17.710	2497	2503	2508	rBV	244016	308716	3.78%	0.534%
45	17.933	2536	2541	2547	rBV	72929	107167	1.31%	0.185%
46	18.010	2550	2554	2566	rVB	22275	42477	0.52%	0.073%
47	18.186	2581	2584	2588	rVB3	9851	10892	0.13%	0.019%
48	18.269	2594	2598	2602	rBV5	7420	11120	0.14%	0.019%
49	18.798	2685	2688	2692	rBV3	10894	15900	0.19%	0.028%
50	18.998	2717	2722	2727	rBV2	48547	66840	0.82%	0.116%
51	19.069	2730	2734	2743	rVV2	45299	66760	0.82%	0.115%
52	19.222	2756	2760	2767	rBV2	73371	108144	1.33%	0.187%
53	19.439	2790	2797	2807	rBV	4760592	6542555	80.18%	11.318%
54	21.269	3102	3108	3118	rBV	2683328	3707028	45.43%	6.413%
55	23.980	3559	3569	3583	rBV2	3369904	8159958	100.00%	14.117%
56	24.174	3593	3602	3619	rVB	1872576	4463420	54.70%	7.722%

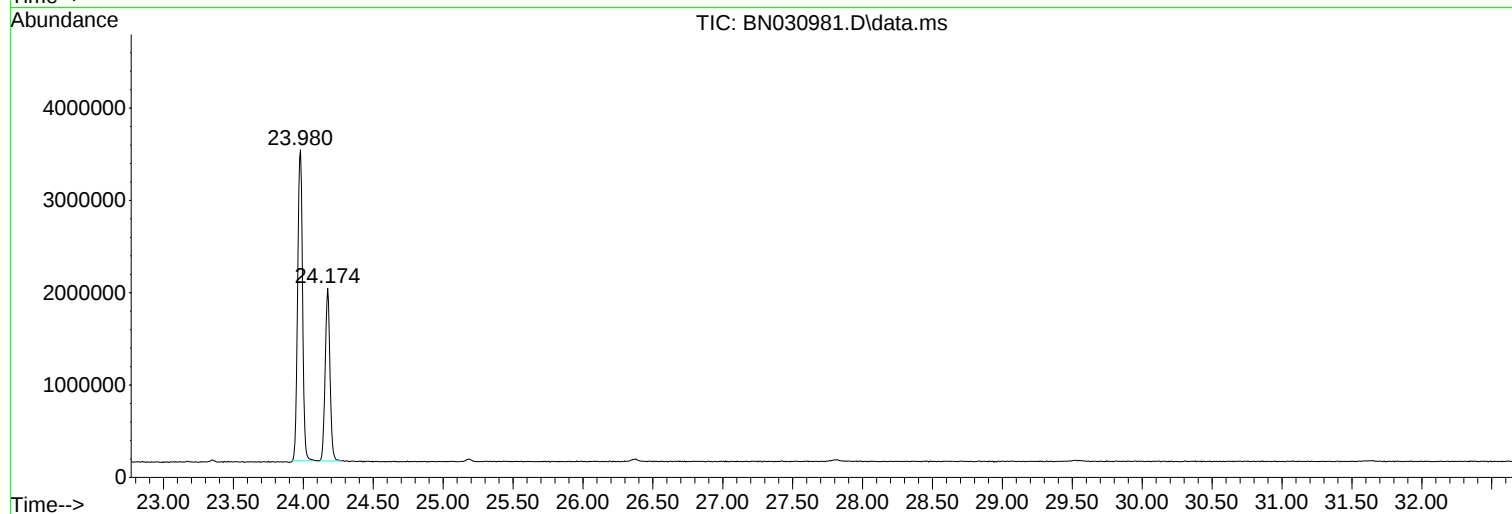
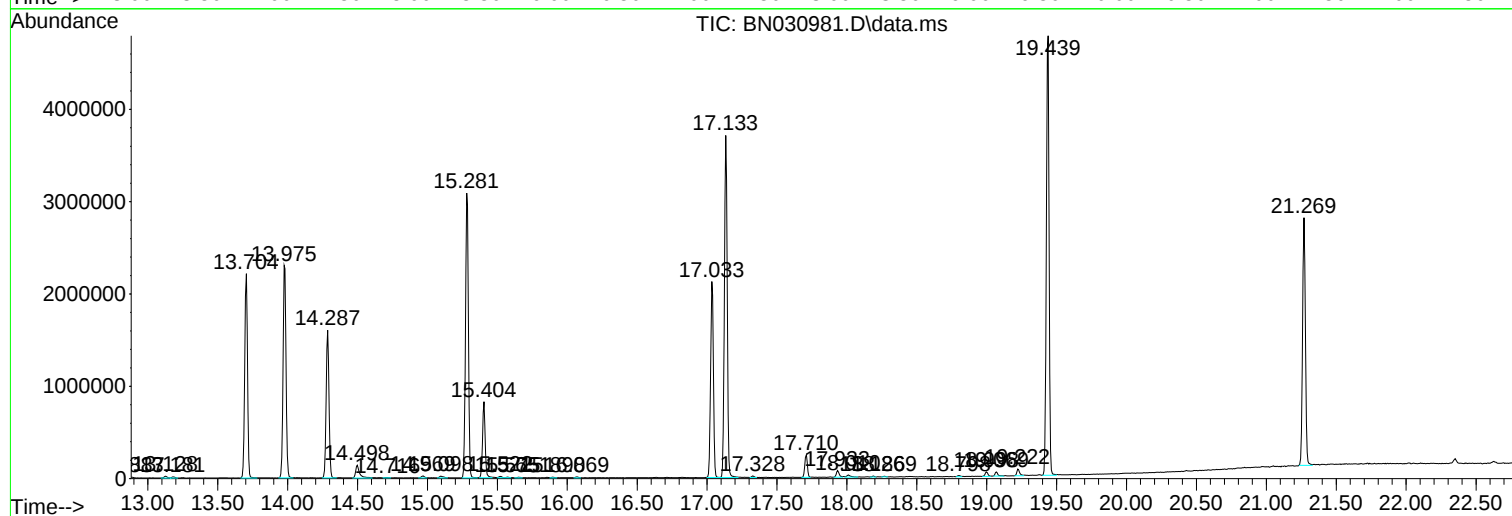
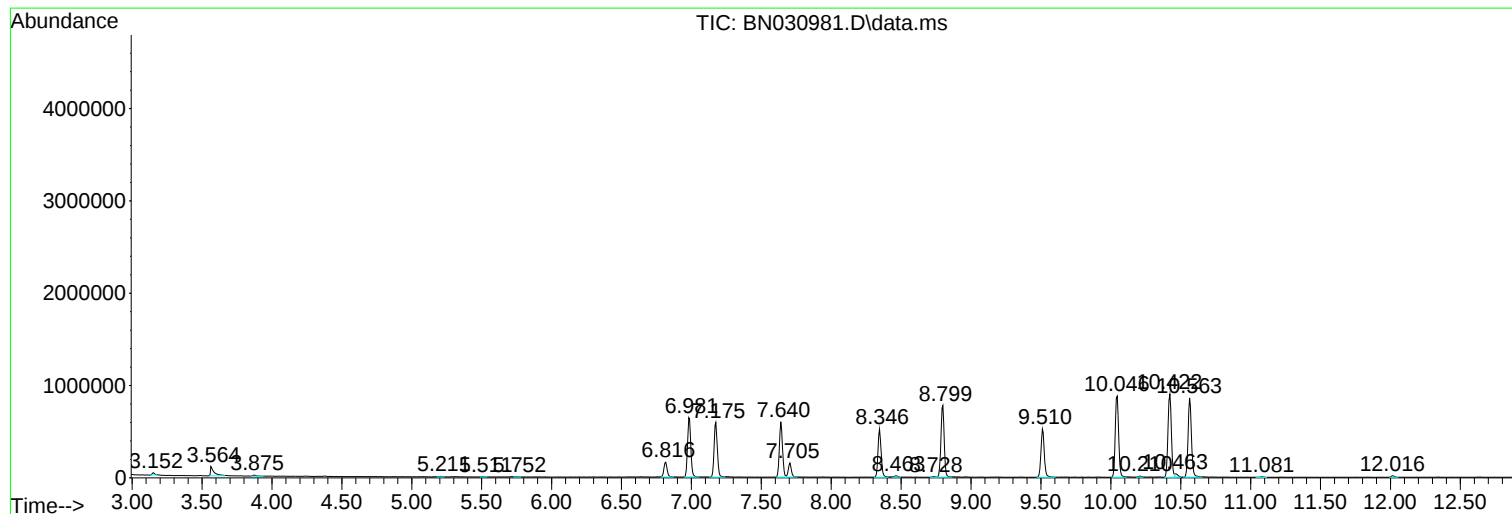
Sum of corrected areas: 57804192

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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



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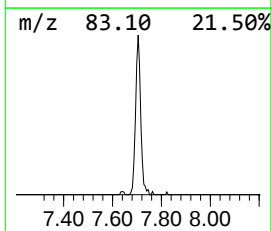
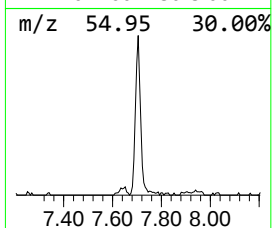
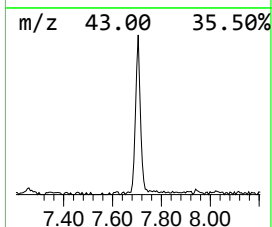
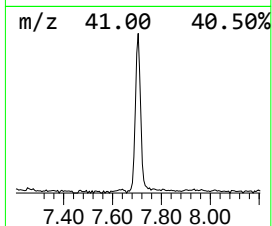
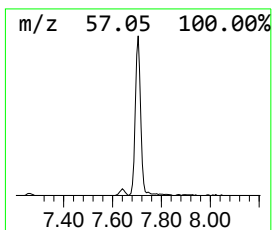
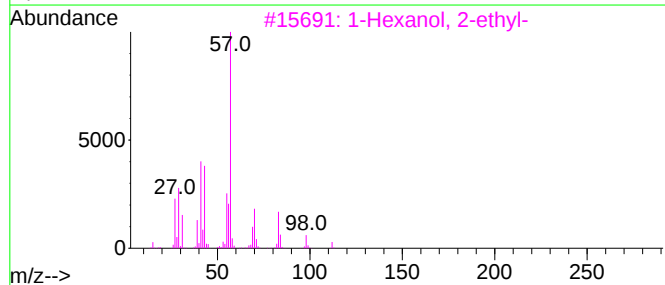
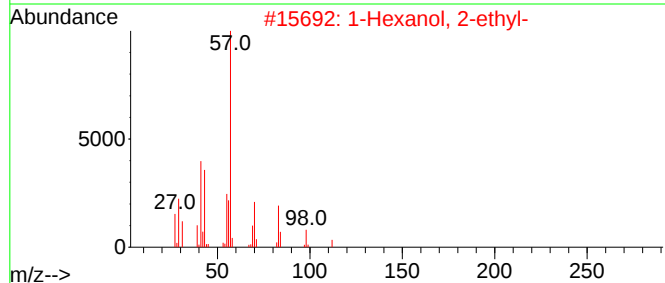
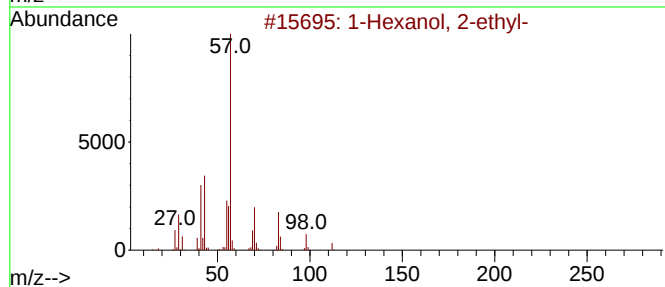
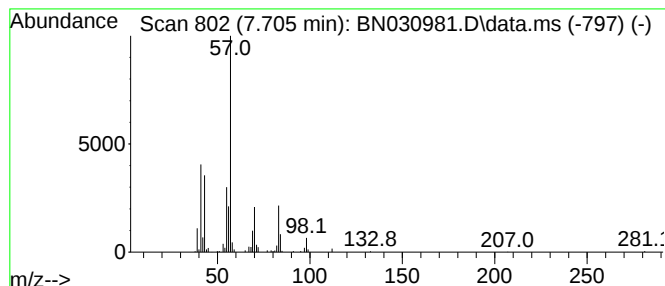
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.705	4.88 ng/ul	227378	1,4-Dichlorobenzene-d4	7.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	90
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
4	Chloroacetic acid, 2-ethylhexyl ...			206	C10H19ClO2	005345-58-4	59
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56



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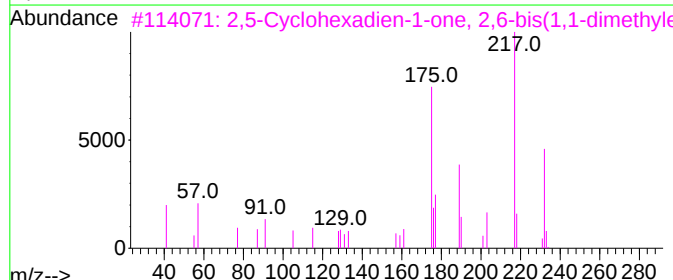
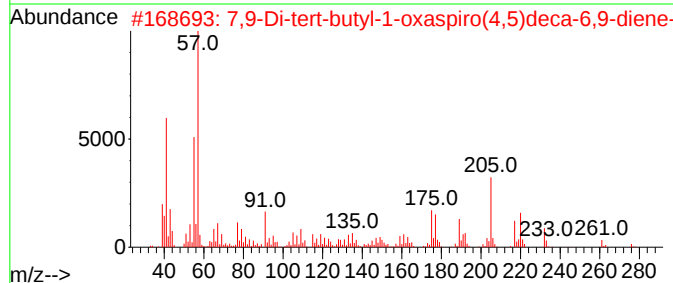
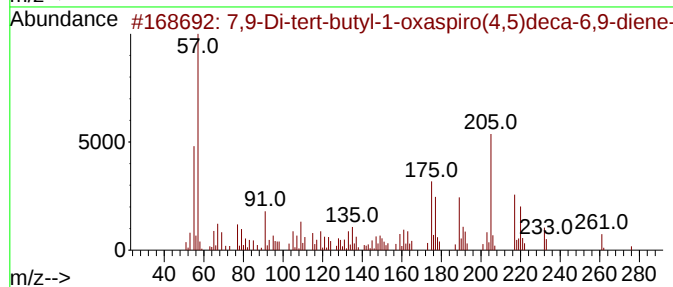
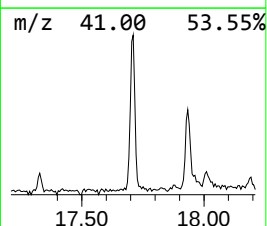
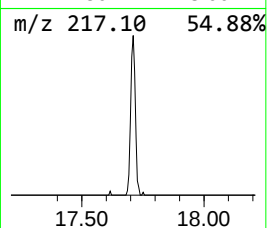
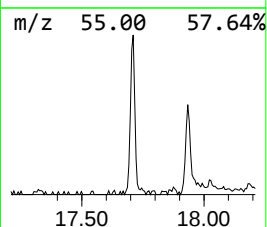
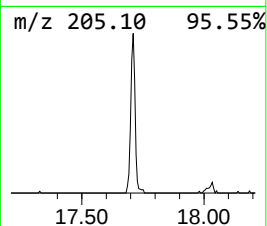
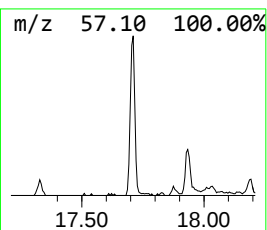
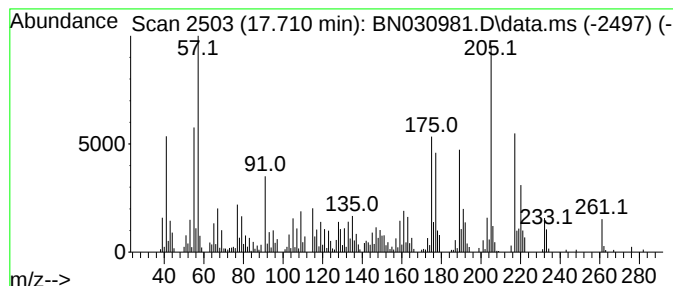
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.710	2.10 ng/ul	308716	Phenanthrene-d10	17.033

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	96		
3	2,5-Cyclohexadien-1-one, 2,6-bis...	232	C16H24O	006738-27-8	83		
4	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	68		
5	9-Acetylphenanthrene	220	C16H12O	002039-77-2	51		



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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...	7.705	4.9	ng/u1	227378	1	7.640	931644	20.0
7,9-Di-tert-but...	17.710	2.1	ng/u1	308716	4	17.033	2944830	20.0