

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014259.D
 Acq On : 23 Mar 2023 15:51
 Operator : CG/JU
 Sample : 01983-02
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYGA9

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-BP030723.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014259.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	34	37	44	rVB	25486	36642	0.56%	0.067%
2	3.858	113	114	132	rBV	45687	94677	1.46%	0.174%
3	4.064	144	149	156	rVB	12271	21975	0.34%	0.040%
4	4.164	161	166	173	rVB	16622	27339	0.42%	0.050%
5	4.275	179	185	188	rBV6	8454	12309	0.19%	0.023%
6	5.099	322	325	333	rVB	11417	21491	0.33%	0.039%
7	6.411	544	548	555	rVB6	8690	13560	0.21%	0.025%
8	7.199	675	682	695	rBV	129153	242414	3.73%	0.445%
9	7.358	702	709	720	rBV	607705	963148	14.81%	1.767%
10	7.563	737	744	754	rBV	542759	908724	13.97%	1.668%
11	8.034	817	824	842	rVB	574620	937062	14.41%	1.720%
12	8.752	937	946	959	rBV	441505	785449	12.08%	1.441%
13	8.881	962	968	973	rVB3	14229	28273	0.43%	0.052%
14	9.210	1015	1024	1039	rBV	783914	1334839	20.52%	2.449%
15	9.593	1081	1089	1096	rVB6	7016	13102	0.20%	0.024%
16	9.934	1139	1147	1162	rBV	625567	1114967	17.14%	2.046%
17	10.481	1233	1240	1262	rBV	792541	1536047	23.62%	2.819%
18	10.857	1296	1304	1315	rBV	879869	1463562	22.50%	2.686%
19	10.999	1320	1328	1350	rBV	581383	1156523	17.78%	2.122%
20	11.546	1416	1421	1427	rBV8	4752	9076	0.14%	0.017%
21	12.446	1569	1574	1583	rBV3	15375	30937	0.48%	0.057%
22	12.563	1589	1594	1600	rBV9	4029	10000	0.15%	0.018%
23	13.310	1713	1721	1730	rVB2	18643	39495	0.61%	0.072%
24	13.522	1751	1757	1761	rBV	41045	60688	0.93%	0.111%
25	13.563	1761	1764	1769	rVB	13015	20303	0.31%	0.037%
26	14.087	1845	1853	1865	rBV	2249874	3042095	46.77%	5.582%
27	14.322	1888	1893	1895	rBV6	5318	7494	0.12%	0.014%
28	14.375	1895	1902	1916	rVV	2262466	3278196	50.40%	6.016%
29	14.681	1946	1954	1962	rBV	1562415	2350696	36.14%	4.314%
30	14.751	1962	1966	1974	rVB	40039	58827	0.90%	0.108%
31	14.916	1987	1994	2010	rBV2	58869	208200	3.20%	0.382%
32	15.351	2063	2068	2076	rBV	26971	46088	0.71%	0.085%
33	15.422	2076	2080	2085	rVV	22844	34869	0.54%	0.064%
34	15.492	2088	2092	2097	rVB	20641	30285	0.47%	0.056%
35	15.675	2116	2123	2136	rBV	3168845	4449649	68.41%	8.165%
36	15.798	2138	2144	2160	rBV	1046859	1555685	23.92%	2.855%

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37	15.916	2161	2164	2168	rVB3	15338	22219	0.34%	0.041%
38	16.039	2179	2185	2192	rVB4	16299	29130	0.45%	0.053%
39	17.022	2347	2352	2357	rBV2	20130	27631	0.42%	0.051%
40	17.116	2364	2368	2372	rBV6	4361	6505	0.10%	0.012%
41	17.228	2384	2387	2392	rBV5	5108	7483	0.12%	0.014%
42	17.386	2410	2414	2416	rBV3	8212	10938	0.17%	0.020%
43	17.439	2416	2423	2433	rVV	2083215	2921669	44.92%	5.361%
44	17.539	2433	2440	2455	rVV2	3708060	5276691	81.13%	9.683%
45	17.710	2465	2469	2475	rVB2	11810	19454	0.30%	0.036%
46	18.086	2528	2533	2539	rBV	609948	749775	11.53%	1.376%
47	18.251	2559	2561	2566	rBV6	10217	10063	0.15%	0.018%
48	18.310	2566	2571	2577	rBV7	25867	52202	0.80%	0.096%
49	18.404	2584	2587	2591	rVB5	17849	23640	0.36%	0.043%
50	18.551	2609	2612	2616	rVB3	21397	23359	0.36%	0.043%
51	18.869	2663	2666	2670	rVB3	22380	26125	0.40%	0.048%
52	19.339	2743	2746	2753	rVB	52658	69992	1.08%	0.128%
53	19.410	2754	2758	2765	rBV	51939	86337	1.33%	0.158%
54	19.533	2776	2779	2782	rBV5	24961	35967	0.55%	0.066%
55	19.763	2812	2818	2830	rBV	4739557	6504344	100.00%	11.936%
56	21.521	3112	3117	3128	rBV	2419362	3361752	51.68%	6.169%
57	23.874	3510	3517	3531	rVB2	3026633	6153011	94.60%	11.291%
58	24.039	3539	3545	3556	rVB2	1457678	3132049	48.15%	5.747%

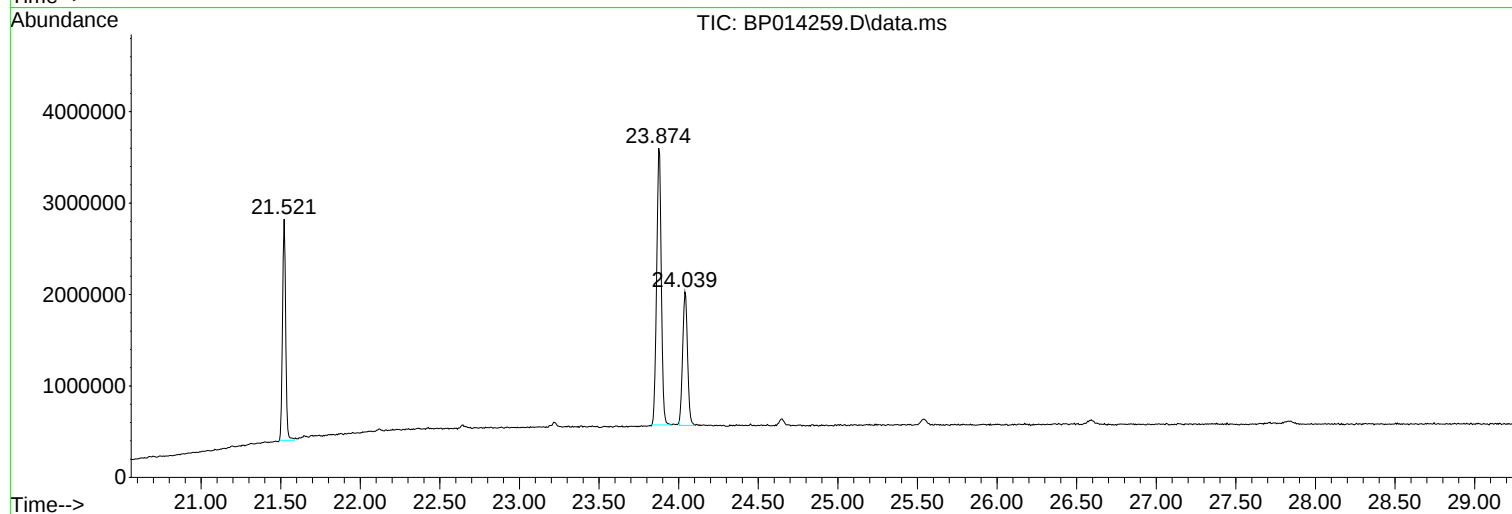
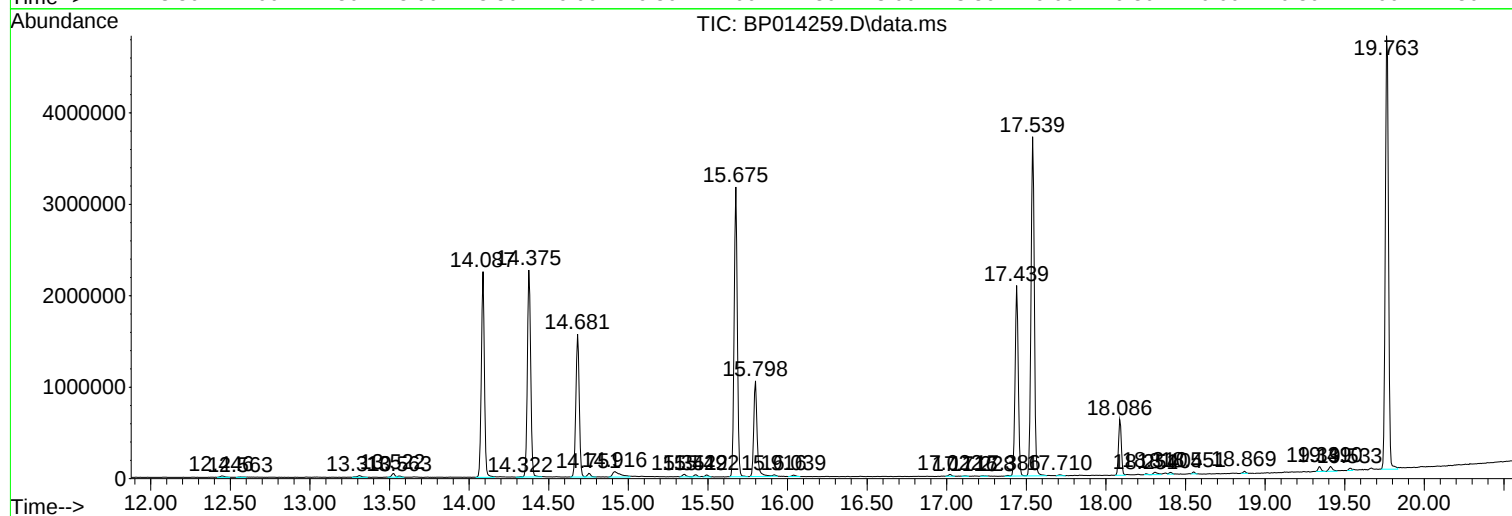
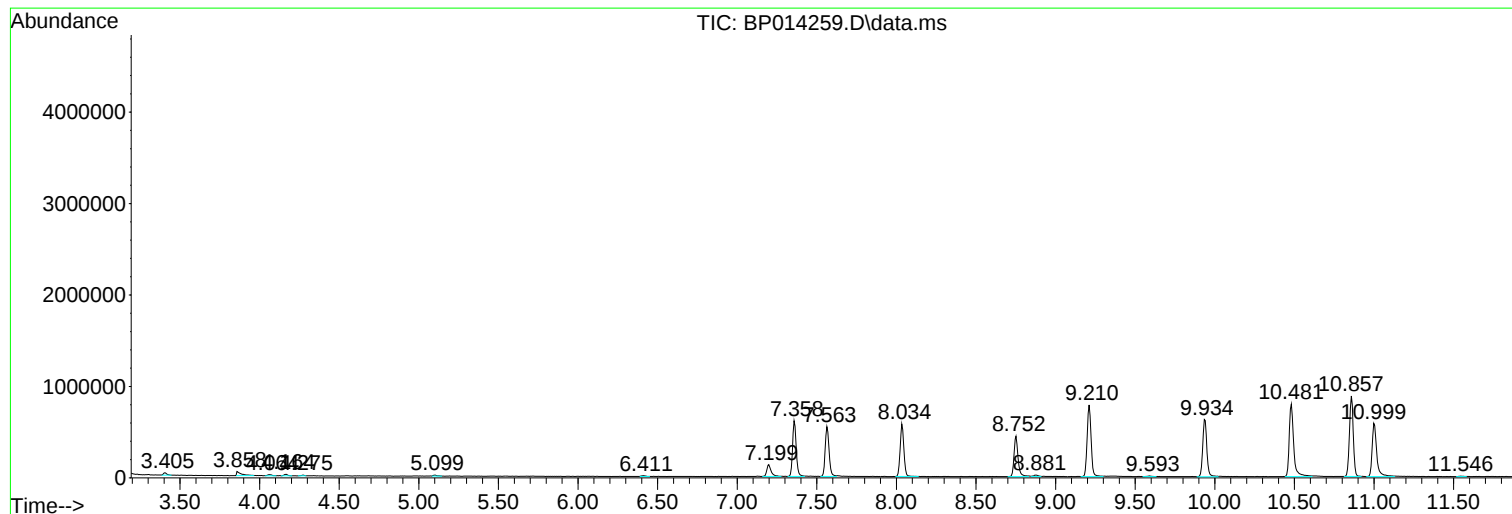
Sum of corrected areas: 54495022

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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



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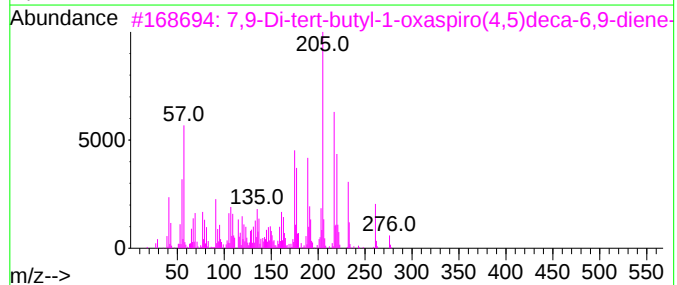
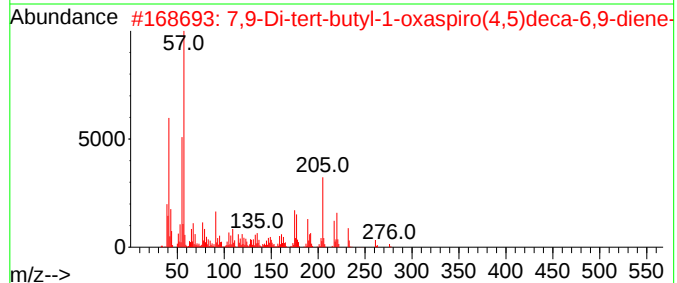
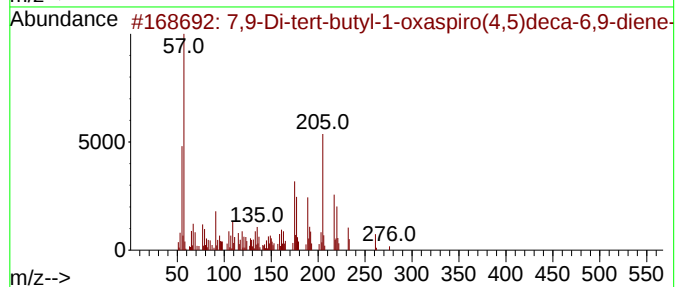
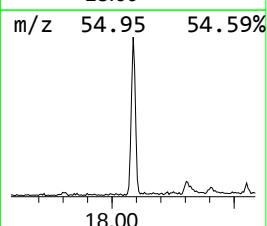
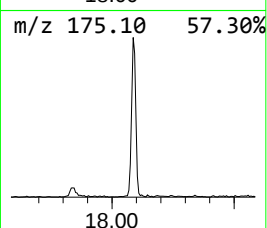
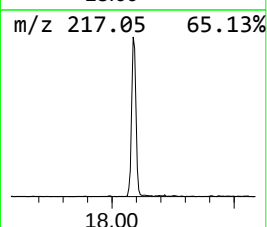
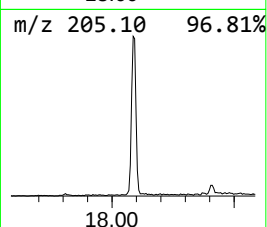
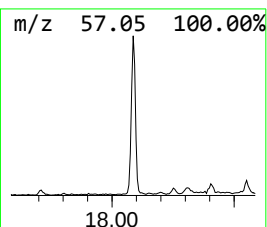
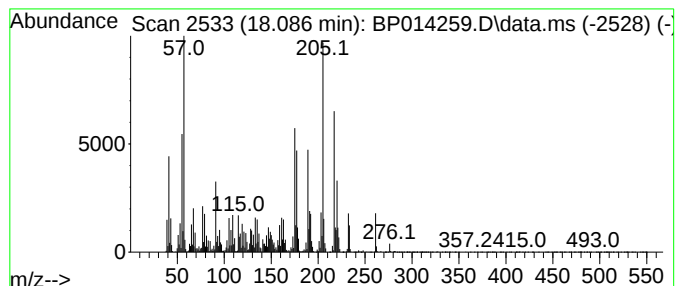
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.086	5.13 ng/ul	749775	Phenanthrene-d10	17.439

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	95		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	92		
4	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	45		
5	Benzenethanol, 0,.beta.,.beta.,2...	220	C15H24O	1010128-94-8	45		



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
7,9-Di-tert-but...	18.086	5.1	ng/ul	749775	4	17.439	2921670	20.0