

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111124\
 Data File : BG063312.D
 Acq On : 12 Nov 2024 6:25
 Operator : RC/JU
 Sample : P4625-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AA5

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_G\Methods\SFAM-EPA-BG110624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063312.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	33	36	42	rVB3	26860	42948	0.77%	0.090%
2	3.704	102	105	113	rBV	105018	157231	2.82%	0.329%
3	4.045	159	163	166	rVB6	8613	11956	0.21%	0.025%
4	4.133	176	178	182	rVB4	8360	9521	0.17%	0.020%
5	4.544	243	248	253	rBV	37830	53851	0.97%	0.113%
6	4.950	316	317	323	rVB3	9420	8823	0.16%	0.018%
7	5.484	405	408	410	rBV4	9637	9690	0.17%	0.020%
8	5.684	438	442	446	rBV3	12571	15896	0.29%	0.033%
9	7.006	661	667	673	rBV	176154	316020	5.68%	0.662%
10	7.165	687	694	705	rVB	450135	738401	13.27%	1.546%
11	7.370	719	729	738	rBV	529292	913667	16.41%	1.913%
12	7.470	742	746	755	rVB5	6904	15991	0.29%	0.033%
13	7.617	765	771	773	rBV2	9482	16527	0.30%	0.035%
14	7.835	800	808	816	rBV	401609	678036	12.18%	1.420%
15	7.899	816	819	822	rVB3	9405	11424	0.21%	0.024%
16	8.540	921	928	938	rBV2	538575	931502	16.73%	1.950%
17	8.927	986	994	998	rBV	79277	138359	2.49%	0.290%
18	8.986	998	1004	1012	rVB	596505	1047491	18.82%	2.193%
19	9.421	1071	1078	1080	rBV5	7535	11084	0.20%	0.023%
20	9.709	1118	1127	1136	rBV2	581891	1007964	18.11%	2.110%
21	10.249	1211	1219	1235	rBV	881826	1542705	27.72%	3.230%
22	10.543	1267	1269	1272	rVB4	10098	8300	0.15%	0.017%
23	10.625	1276	1283	1290	rBV	633333	1117295	20.07%	2.339%
24	10.761	1298	1306	1315	rBV	668273	1257084	22.58%	2.632%
25	11.795	1480	1482	1488	rVB4	5715	8615	0.15%	0.018%
26	12.212	1549	1553	1555	rBV3	11473	14579	0.26%	0.031%
27	12.735	1639	1642	1646	rBV4	9233	14380	0.26%	0.030%
28	12.805	1651	1654	1655	rBV2	7998	7913	0.14%	0.017%
29	12.823	1655	1657	1662	rVB3	7051	12213	0.22%	0.026%
30	13.075	1696	1700	1708	rBV5	12444	23085	0.41%	0.048%
31	13.393	1749	1754	1759	rBV3	23118	42470	0.76%	0.089%
32	13.804	1821	1824	1827	rBV2	5739	8053	0.14%	0.017%
33	13.880	1827	1837	1848	rVV	1778981	2598886	46.69%	5.441%
34	14.162	1878	1885	1893	rBV	1885287	2903229	52.16%	6.078%
35	14.474	1929	1938	1947	rBV	1284880	1968667	35.37%	4.122%
36	14.556	1947	1952	1957	rVB4	21936	33566	0.60%	0.070%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Title : SVOA CALIBRATION

37	14.685	1969	1974	1984	rBV2	124625	245464	4.41%	0.514%
38	15.144	2050	2052	2058	rVB4	7280	12008	0.22%	0.025%
39	15.273	2073	2074	2081	rVB5	8733	15315	0.28%	0.032%
40	15.467	2100	2107	2114	rBV	2582874	3846831	69.11%	8.054%
41	15.590	2122	2128	2138	rBV	777114	1194055	21.45%	2.500%
42	15.708	2145	2148	2152	rVB5	12935	16361	0.29%	0.034%
43	15.837	2166	2170	2173	rVB6	9398	11362	0.20%	0.024%
44	16.172	2222	2227	2229	rBV5	7831	12585	0.23%	0.026%
45	16.313	2248	2251	2257	rBV3	5315	9173	0.16%	0.019%
46	16.360	2257	2259	2264	rVB4	7345	9544	0.17%	0.020%
47	16.824	2333	2338	2341	rVB4	14447	22665	0.41%	0.047%
48	16.900	2349	2351	2356	rVB4	8428	8599	0.15%	0.018%
49	16.977	2363	2364	2368	rBV3	7601	8313	0.15%	0.017%
50	17.147	2389	2393	2395	rBV3	7348	9372	0.17%	0.020%
51	17.177	2395	2398	2399	rVV2	11029	12267	0.22%	0.026%
52	17.224	2400	2406	2414	rVV	1691744	2461655	44.22%	5.154%
53	17.323	2416	2423	2434	rVB	2991279	4559174	81.91%	9.545%
54	17.506	2451	2454	2458	rVB2	6752	8417	0.15%	0.018%
55	17.611	2465	2472	2477	rBV9	6923	13618	0.24%	0.029%
56	17.893	2513	2520	2526	rBV2	365149	462933	8.32%	0.969%
57	18.193	2567	2571	2574	rBV	14502	18136	0.33%	0.038%
58	18.622	2638	2644	2646	rBV5	8700	9892	0.18%	0.021%
59	19.186	2735	2740	2747	rVB3	48037	75066	1.35%	0.157%
60	19.256	2747	2752	2759	rBV2	42663	75255	1.35%	0.158%
61	19.433	2780	2782	2785	rBV4	6804	7598	0.14%	0.016%
62	19.538	2796	2800	2801	rBV4	11168	14441	0.26%	0.030%
63	19.621	2808	2814	2826	rBV	3875536	5566293	100.00%	11.654%
64	19.850	2850	2853	2861	rVB8	29953	49546	0.89%	0.104%
65	20.273	2922	2925	2927	rBV4	15729	17115	0.31%	0.036%
66	21.477	3124	3130	3144	rVB2	1853813	2868880	51.54%	6.007%
67	24.309	3600	3612	3625	rBV	2122304	5524140	99.24%	11.566%
68	24.515	3638	3647	3659	rVB	1091441	2899126	52.08%	6.070%

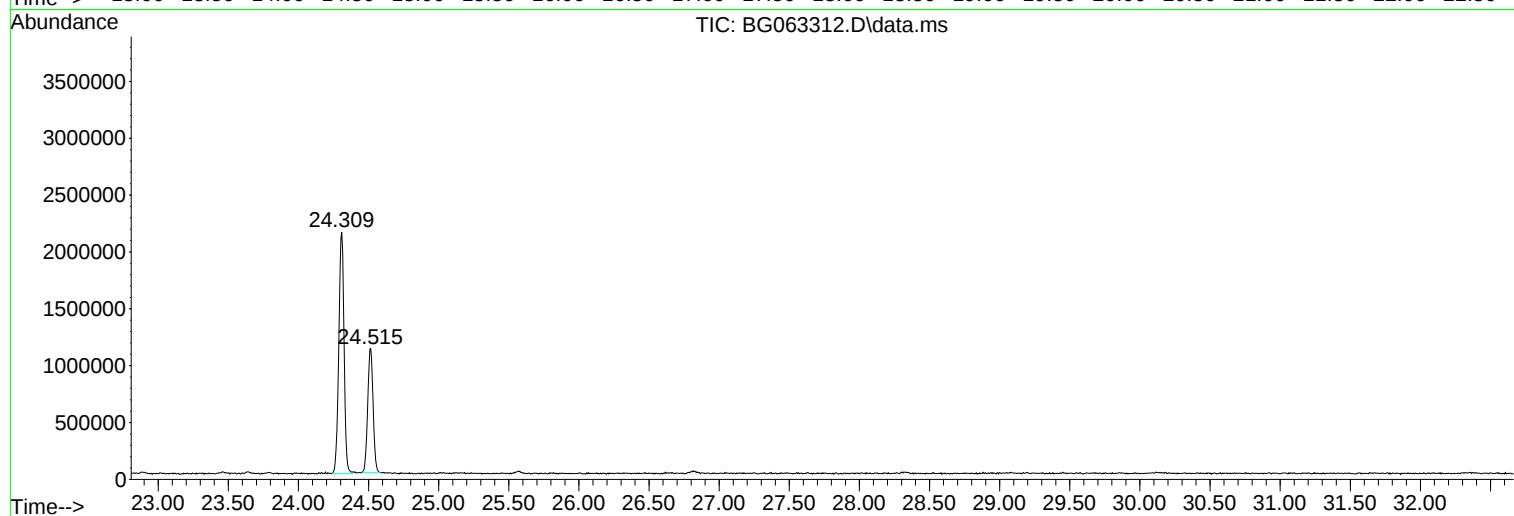
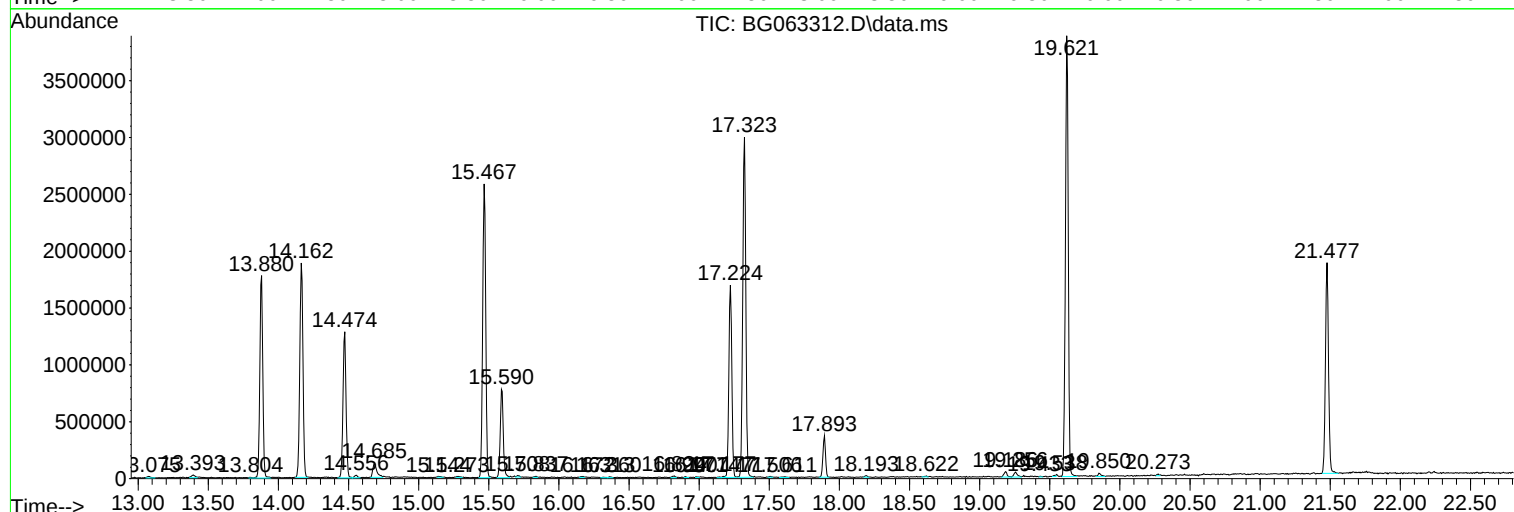
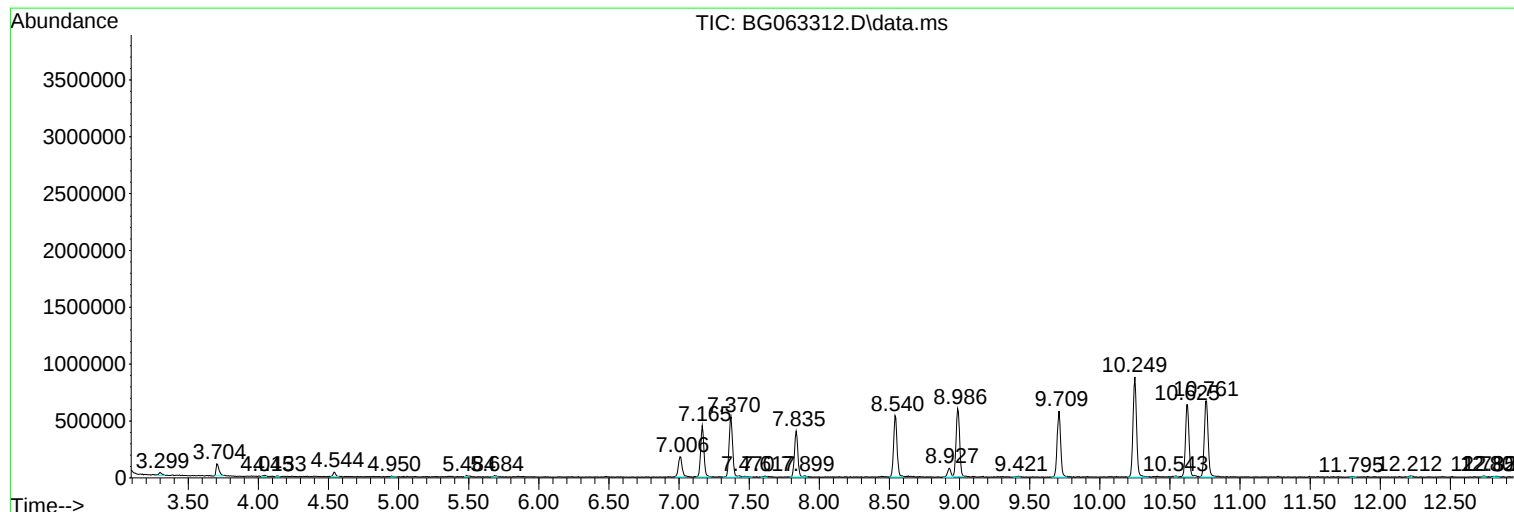
Sum of corrected areas: 47762621

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BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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



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ClientSampleId :
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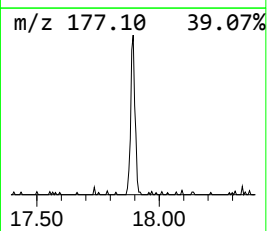
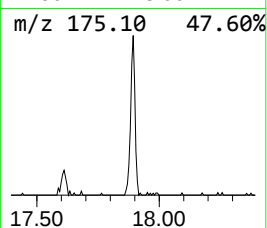
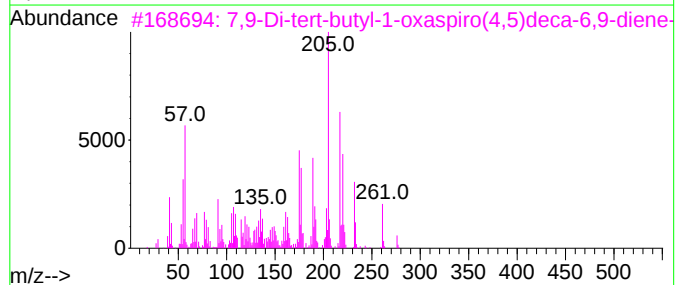
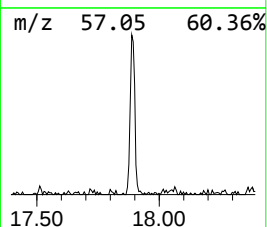
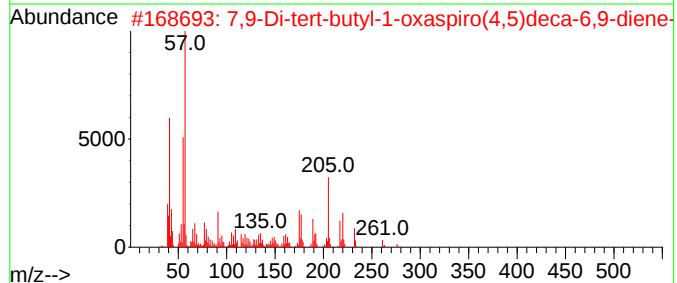
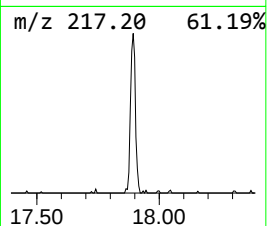
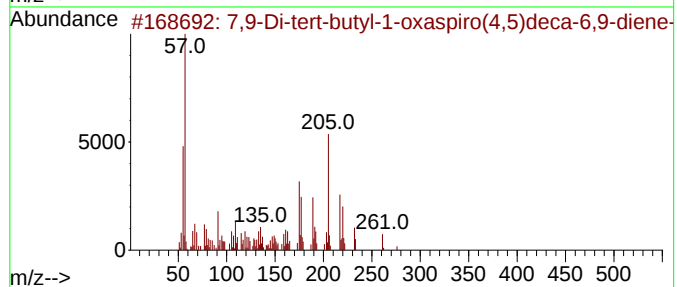
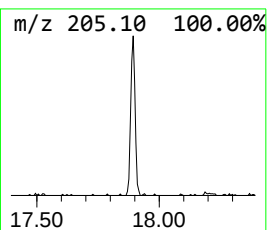
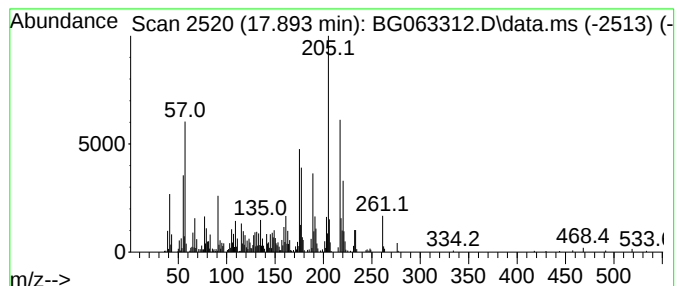
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.893	3.76 ng/ul	462933	Phenanthrene-d10	17.224

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	94		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	86		
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	83		
4	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	70		
5	2H-1-Benzopyran, 6,7-dimethoxy-2...	220	C13H16O3	000644-06-4	55		



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
7,9-Di-tert-but...	17.893	3.8	ng/ul	462933	4	17.224	2461660	20.0