

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122024\
 Data File : BG063772.D
 Acq On : 20 Dec 2024 15:20
 Operator : RC/JU
 Sample : P5305-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E2AX4

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-BG121124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063772.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	27	30	38	rVB2	37026	61588	1.29%	0.150%
2	3.669	95	99	108	rBV	139582	231065	4.83%	0.563%
3	4.686	270	272	275	rVB3	8895	9167	0.19%	0.022%
4	6.190	525	528	532	rBV3	11233	13860	0.29%	0.034%
5	6.989	658	664	673	rVB	175394	325745	6.81%	0.794%
6	7.130	683	688	697	rVB	393662	648149	13.55%	1.580%
7	7.336	718	723	732	rVB	416079	704823	14.73%	1.719%
8	7.800	796	802	811	rVB	436288	705068	14.74%	1.719%
9	8.370	897	899	902	rVB2	6963	7158	0.15%	0.017%
10	8.517	918	924	936	rBV2	433371	756630	15.82%	1.845%
11	8.957	992	999	1008	rBV	482916	816359	17.06%	1.991%
12	9.380	1067	1071	1072	rBV3	6101	7248	0.15%	0.018%
13	9.680	1115	1122	1130	rBV	358713	637085	13.32%	1.553%
14	10.050	1183	1185	1190	rVB4	5973	8392	0.18%	0.020%
15	10.091	1190	1192	1195	rBV2	6241	6484	0.14%	0.016%
16	10.226	1208	1215	1223	rBV	534532	987591	20.64%	2.408%
17	10.591	1270	1277	1287	rVB	589900	1035762	21.65%	2.526%
18	10.726	1293	1300	1311	rBV	439100	843693	17.64%	2.057%
19	11.260	1388	1391	1394	rVB3	6392	7549	0.16%	0.018%
20	11.425	1416	1419	1421	rBV	8053	9294	0.19%	0.023%
21	11.554	1439	1441	1443	rBV2	6788	7010	0.15%	0.017%
22	12.900	1667	1670	1673	rVB2	6277	9596	0.20%	0.023%
23	12.970	1680	1682	1686	rBV4	5166	8059	0.17%	0.020%
24	13.264	1727	1732	1734	rBV4	4592	8005	0.17%	0.020%
25	13.417	1756	1758	1762	rVB4	6742	7865	0.16%	0.019%
26	13.581	1783	1786	1788	rBV2	5404	6524	0.14%	0.016%
27	13.622	1790	1793	1796	rVB3	6122	7542	0.16%	0.018%
28	13.846	1825	1831	1839	rBV	1135465	1696280	35.46%	4.136%
29	14.134	1873	1880	1888	rBV	1326313	2070899	43.29%	5.050%
30	14.439	1923	1932	1941	rBV	1113078	1687831	35.28%	4.116%
31	14.680	1968	1973	1986	rBV2	104727	250876	5.24%	0.612%
32	15.373	2087	2091	2095	rBV2	4494	7588	0.16%	0.019%
33	15.438	2095	2102	2111	rBV	1995245	2983716	62.37%	7.276%
34	15.573	2118	2125	2135	rBV	535343	880460	18.40%	2.147%
35	15.802	2159	2164	2170	rBV2	43060	60622	1.27%	0.148%
36	16.977	2358	2364	2367	rBV3	9330	16544	0.35%	0.040%

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Peak Location: TOP

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Title : SVOA CALIBRATION

37	17.189	2391	2400	2409	rBV2	1577466	2489790	52.05%	6.071%
38	17.295	2411	2418	2428	rVB	2564676	3988311	83.37%	9.725%
39	17.865	2509	2515	2523	rVB5	12370	27596	0.58%	0.067%
40	18.387	2601	2604	2608	rVB6	8414	11840	0.25%	0.029%
41	18.517	2622	2626	2629	rBV4	8522	9665	0.20%	0.024%
42	19.151	2730	2734	2738	rBV3	41517	56755	1.19%	0.138%
43	19.222	2742	2746	2752	rVB3	35116	50865	1.06%	0.124%
44	19.516	2792	2796	2799	rVB6	13408	17027	0.36%	0.042%
45	19.592	2802	2809	2821	rBV	3360018	4783825	100.00%	11.665%
46	19.815	2844	2847	2848	rBV3	6695	6704	0.14%	0.016%
47	20.561	2971	2974	2977	rBV5	28169	30782	0.64%	0.075%
48	21.437	3117	3123	3139	rVB2	1702817	2762731	57.75%	6.737%
49	22.812	3351	3357	3371	rBV	909038	1967147	41.12%	4.797%
50	24.251	3592	3602	3614	rBV	1693318	4381613	91.59%	10.684%
51	24.451	3628	3636	3651	rVB	1056274	2893209	60.48%	7.055%

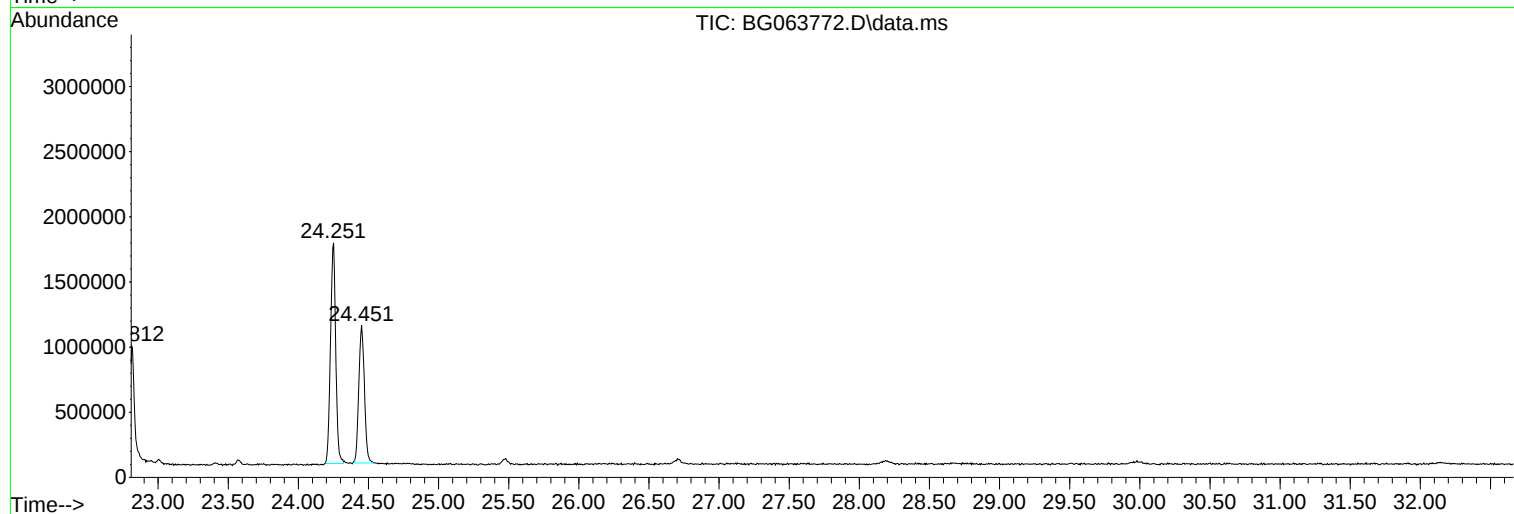
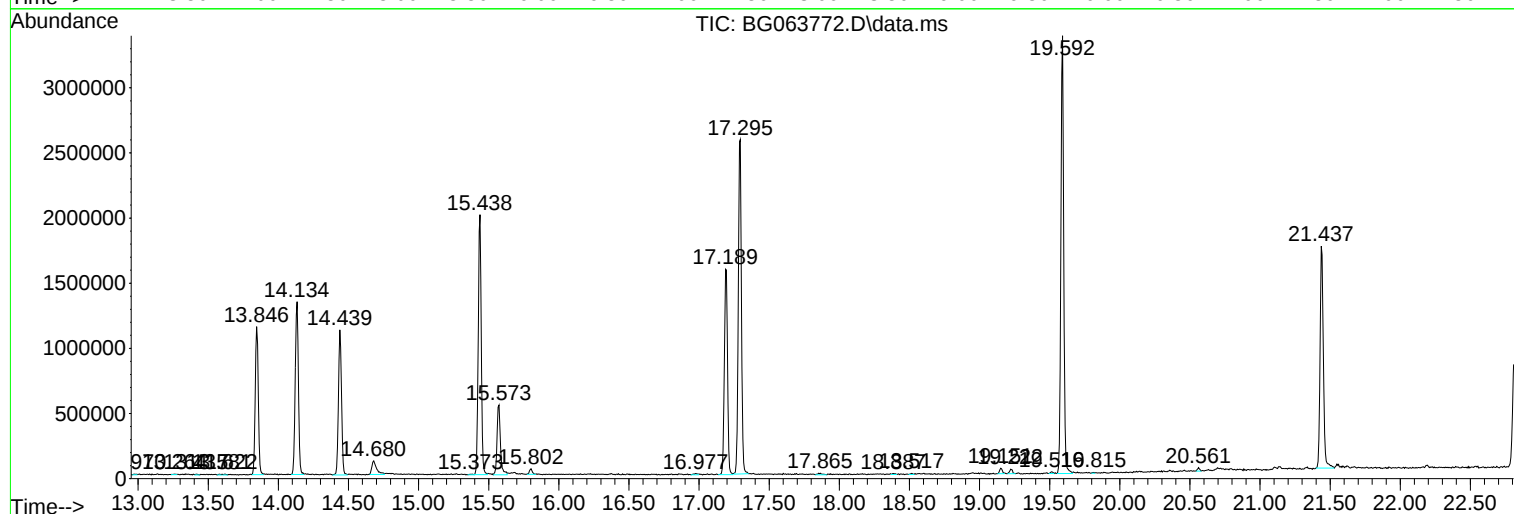
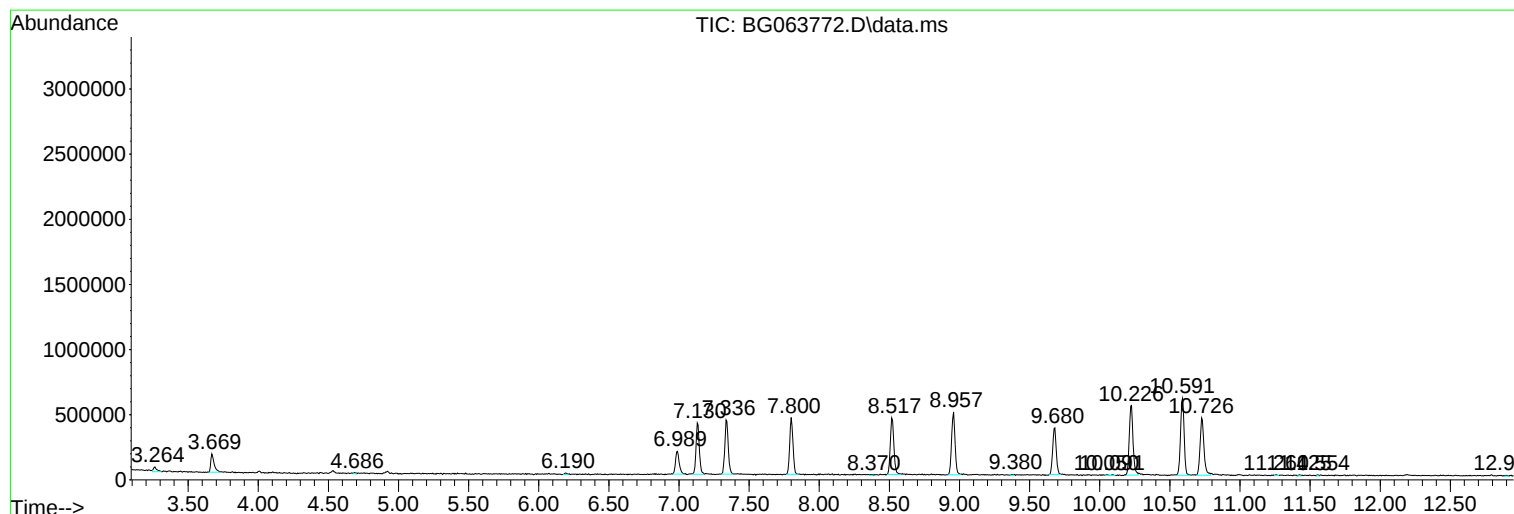
Sum of corrected areas: 41009987

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION

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



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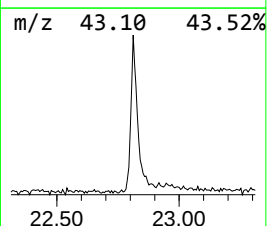
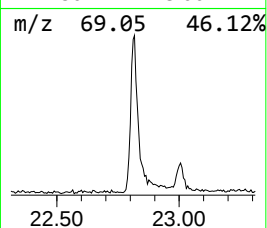
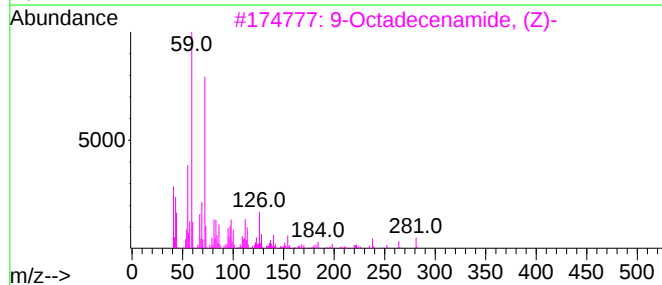
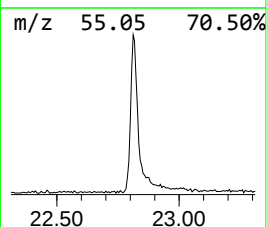
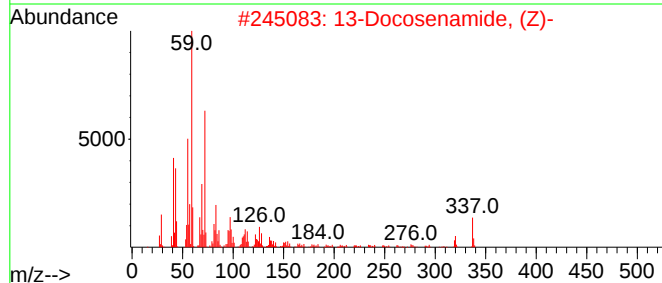
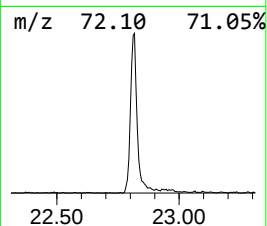
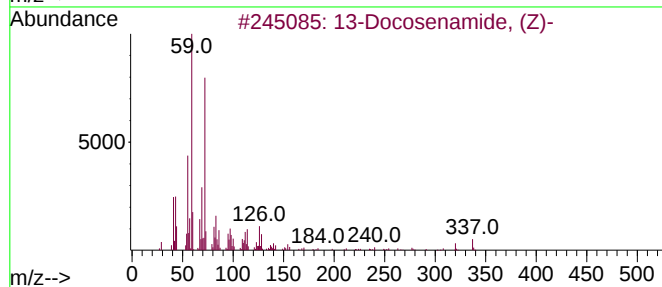
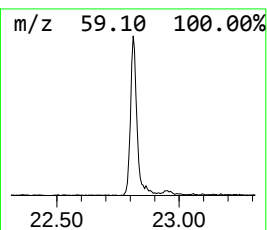
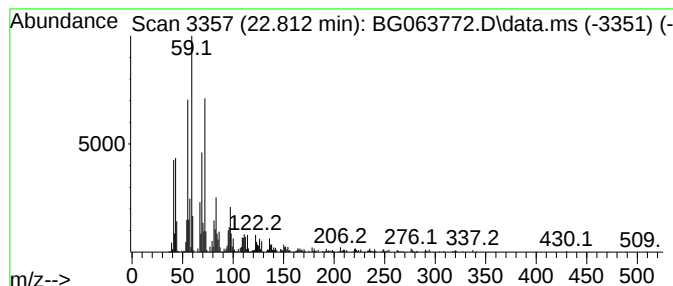
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.812	14.24 ng/ul	1967150	Chrysene-d12	21.437

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	95
2			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	62
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	53
4			2-Octanol, 2-methyl-6-methylene-	156	C10H20O	018479-59-9	46
5			Undecanol-3	172	C11H24O	006929-08-4	35



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
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
13-Docosenamide...	22.812	14.2	ng/ul	1967150	5	21.437	2762730	20.0