

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060324\
 Data File : BM045804.D
 Acq On : 03 Jun 2024 14:29
 Operator : MA/JU
 Sample : P2663-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BHBZ5

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_M\Methods\SFAM-EPA-BM053124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM045804.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.057	9	12	19	rVB	54661	72735	1.66%	0.165%
2	3.493	83	86	105	rBV	136949	366734	8.38%	0.831%
3	3.746	126	129	134	rVB	27641	34836	0.80%	0.079%
4	4.175	198	202	207	rVB	18994	25112	0.57%	0.057%
5	4.987	337	340	343	rVB3	10237	12778	0.29%	0.029%
6	5.763	468	472	477	rVB6	11571	17397	0.40%	0.039%
7	5.957	504	505	508	rBV3	4173	4551	0.10%	0.010%
8	6.240	549	553	561	rVB8	9614	22773	0.52%	0.052%
9	6.628	611	619	620	rBV7	4362	9483	0.22%	0.021%
10	6.739	631	638	646	rBV2	110776	243598	5.57%	0.552%
11	6.834	649	654	671	rVB	783537	1398974	31.96%	3.169%
12	7.057	684	692	712	rBV2	212721	766589	17.51%	1.737%
13	7.469	755	762	774	rVB	895213	1392130	31.80%	3.154%
14	7.975	843	848	850	rBV6	4112	5363	0.12%	0.012%
15	8.251	888	895	911	rBV	390215	745662	17.03%	1.689%
16	8.581	948	951	952	rBV3	7055	7708	0.18%	0.017%
17	8.633	953	960	975	rVV	994414	1730844	39.54%	3.921%
18	8.757	975	981	990	rVV4	20628	58670	1.34%	0.133%
19	8.833	991	994	1001	rVV7	14694	30253	0.69%	0.069%
20	9.245	1060	1064	1069	rBV3	14184	29050	0.66%	0.066%
21	9.363	1080	1084	1087	rBV6	6200	9390	0.21%	0.021%
22	9.916	1170	1178	1194	rBV2	261018	620352	14.17%	1.405%
23	10.239	1225	1233	1243	rBV	1093923	1798633	41.09%	4.075%
24	10.422	1257	1264	1283	rBV	854547	1738685	39.72%	3.939%
25	11.439	1435	1437	1442	rVB5	3216	4695	0.11%	0.011%
26	11.663	1471	1475	1476	rBV4	6464	5922	0.14%	0.013%
27	11.745	1481	1489	1490	rBV3	14421	28151	0.64%	0.064%
28	11.839	1501	1505	1511	rVB3	20373	34837	0.80%	0.079%
29	12.669	1643	1646	1650	rBV6	3042	4591	0.10%	0.010%
30	13.086	1710	1717	1723	rBV10	9081	21632	0.49%	0.049%
31	13.439	1774	1777	1781	rVB5	5506	7275	0.17%	0.016%
32	13.563	1792	1798	1811	rBV	1388839	2002544	45.75%	4.537%
33	13.745	1821	1829	1832	rBV5	10307	21960	0.50%	0.050%
34	13.804	1832	1839	1848	rVV	2156072	3036616	69.37%	6.879%
35	14.110	1885	1891	1899	rBV	1511479	2149997	49.12%	4.871%
36	14.180	1899	1903	1911	rVB	68524	94919	2.17%	0.215%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Title : SVOA CALIBRATION

37	14.298	1918	1923	1927	rVB7	3558	5262	0.12%	0.012%
38	14.439	1942	1947	1949	rBV6	3765	5617	0.13%	0.013%
39	14.974	2032	2038	2041	rBV7	4939	7550	0.17%	0.017%
40	15.110	2053	2061	2074	rBV	2595944	3616284	82.61%	8.192%
41	15.357	2098	2103	2107	rVB6	11009	17238	0.39%	0.039%
42	16.356	2268	2273	2275	rBV5	3723	6138	0.14%	0.014%
43	16.539	2300	2304	2308	rVB7	5633	7860	0.18%	0.018%
44	16.651	2320	2323	2325	rBV4	5234	5091	0.12%	0.012%
45	16.856	2352	2358	2367	rBV	1833262	2323041	53.07%	5.263%
46	16.956	2369	2375	2390	rVV	3022861	4078452	93.17%	9.239%
47	17.180	2410	2413	2416	rBV5	4502	5975	0.14%	0.014%
48	17.562	2475	2478	2481	rBV5	5055	8086	0.18%	0.018%
49	17.721	2501	2505	2507	rBV5	3065	5192	0.12%	0.012%
50	18.821	2688	2692	2698	rVB3	30978	45685	1.04%	0.103%
51	18.892	2700	2704	2708	rVB	34299	43146	0.99%	0.098%
52	19.215	2757	2759	2760	rBV2	8614	6517	0.15%	0.015%
53	19.256	2760	2766	2775	rVB	3240418	4377308	100.00%	9.917%
54	19.503	2806	2808	2813	rVB5	11839	14063	0.32%	0.032%
55	20.821	3028	3032	3040	rBV	176456	218706	5.00%	0.495%
56	21.068	3069	3074	3080	rBV	1623838	2148578	49.08%	4.867%
57	22.356	3287	3293	3308	rBV	1343598	2402326	54.88%	5.442%
58	23.609	3498	3506	3518	rBV2	1739976	3810844	87.06%	8.633%
59	23.791	3529	3537	3547	rBV	1094385	2427086	55.45%	5.498%

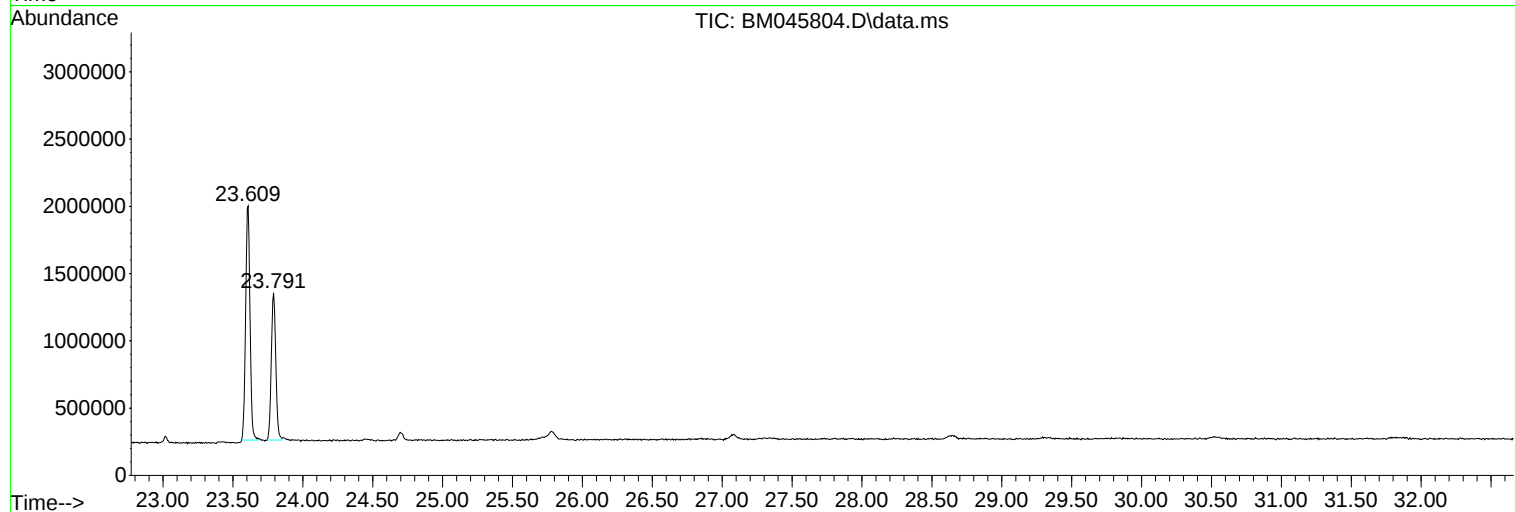
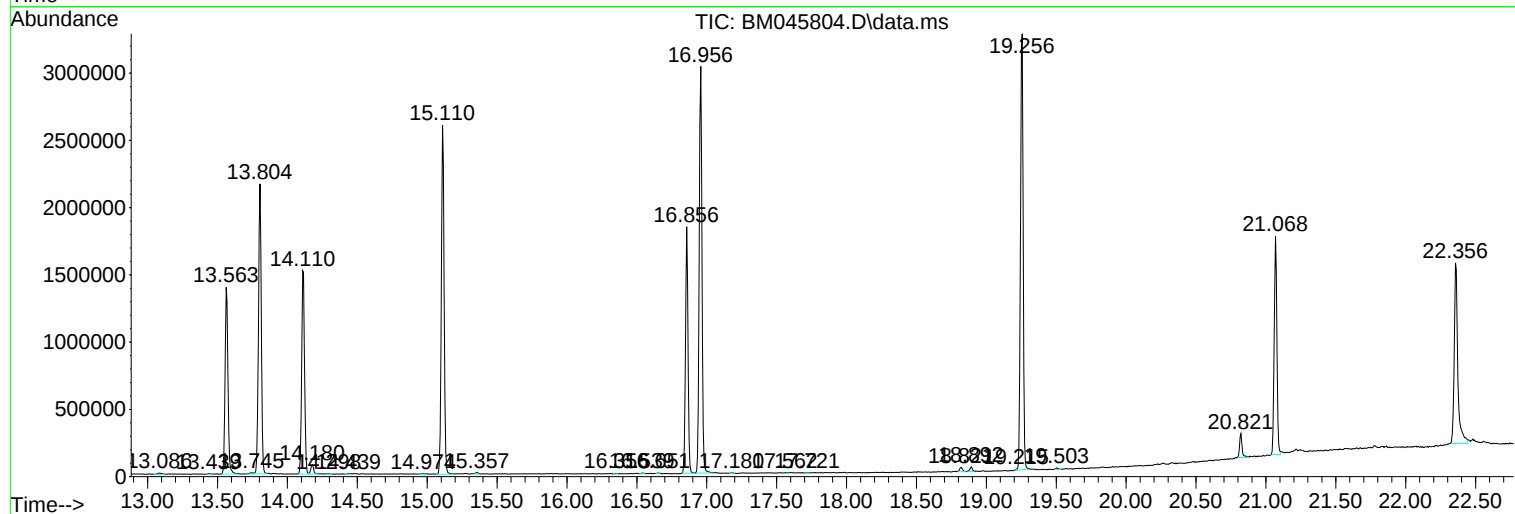
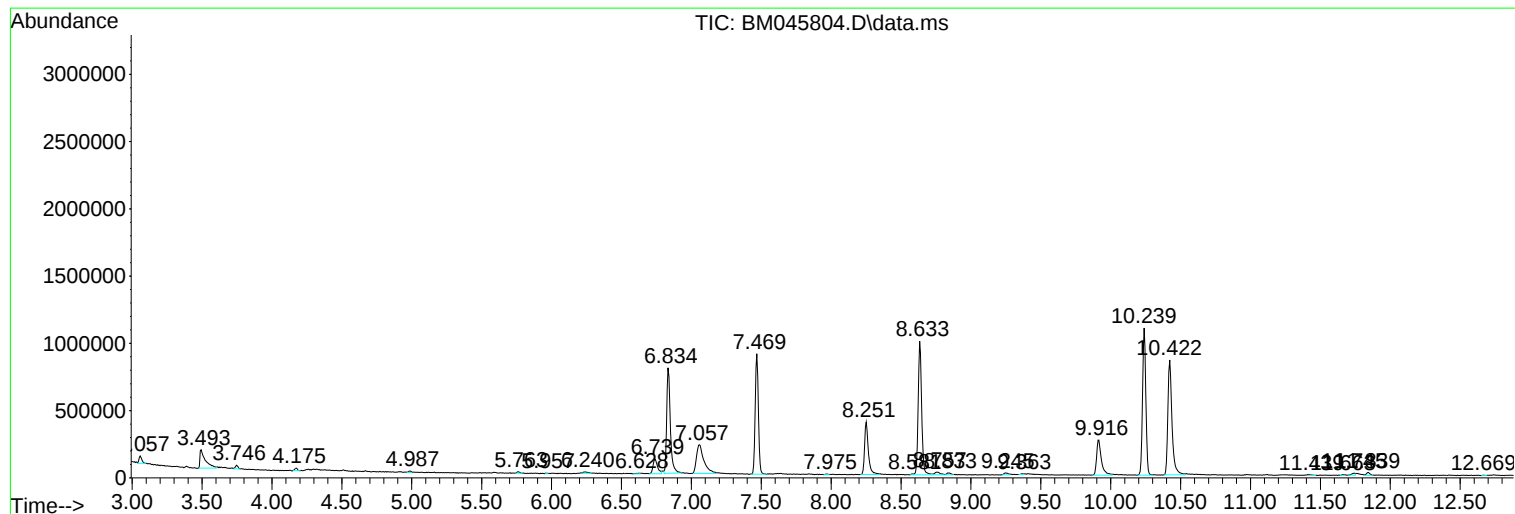
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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



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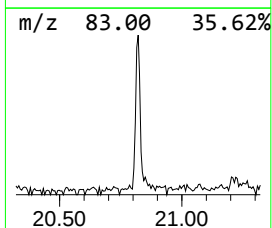
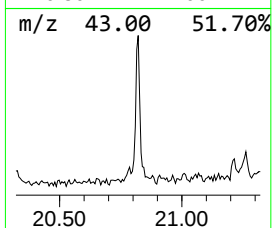
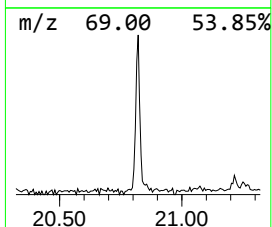
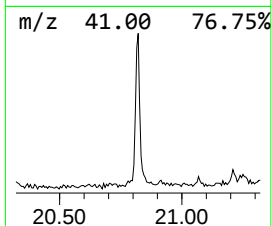
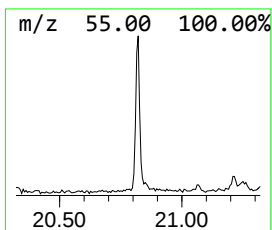
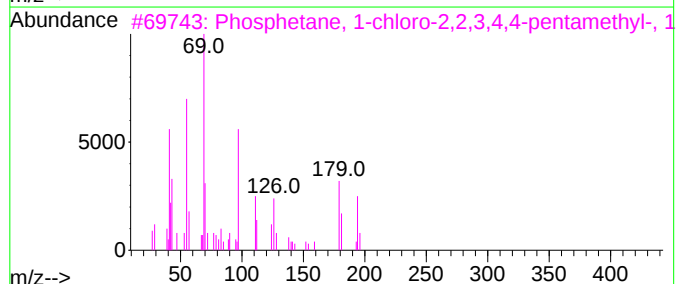
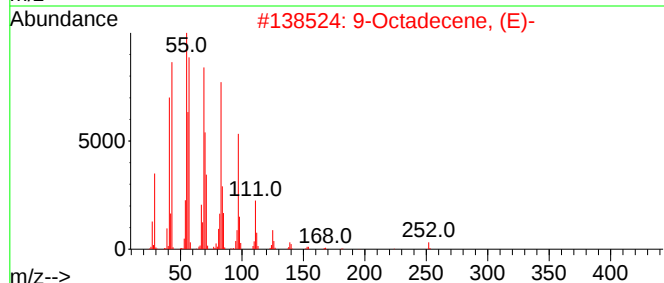
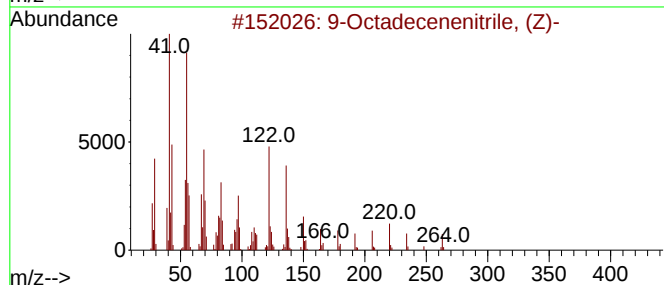
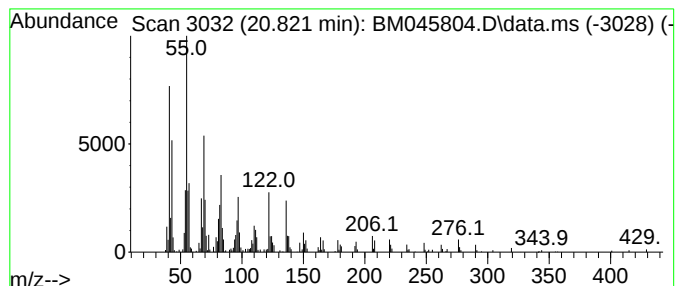
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 9-Octadecenitrile, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.821	2.04 ng/ul	218706	Chrysene-d12	21.068

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenitrile, (Z)-	263	C18H33N	000112-91-4	97
2			9-Octadecene, (E)-	252	C18H36	007206-25-9	25
3			Phosphetane, 1-chloro-2,2,3,4,4-...	194	C8H16ClOP	029276-11-7	15
4			2-Pentenal, 2-methyl-4-oxo-5-(2,...	262	C15H18O4	055721-95-4	15
5			Cyclohexane, 1-(cyclohexylmethyl...)	208	C15H28	054934-92-8	15



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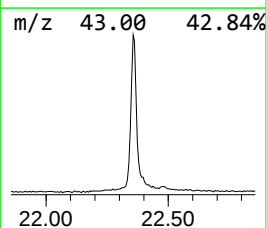
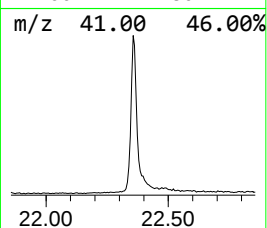
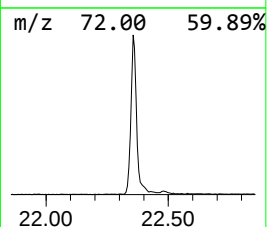
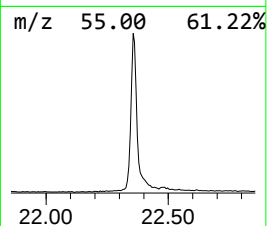
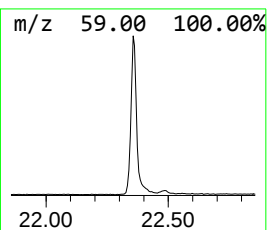
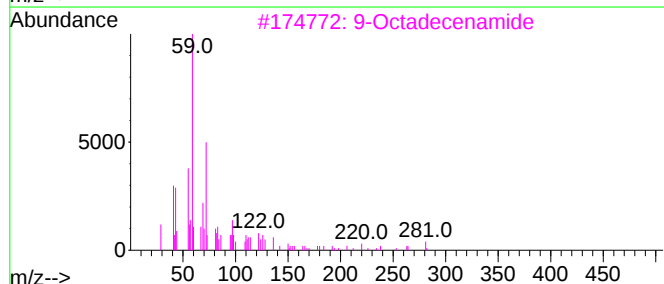
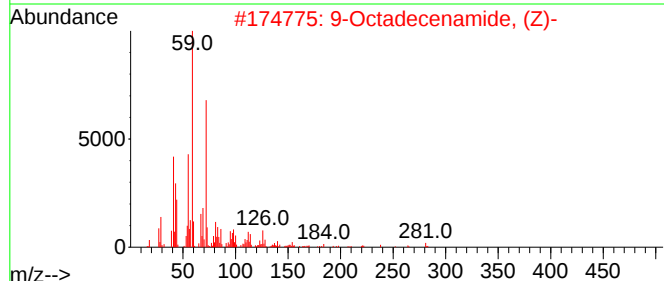
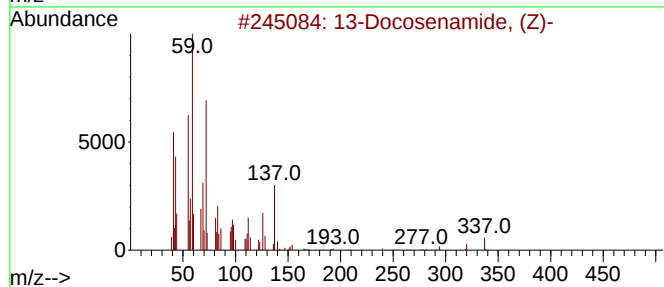
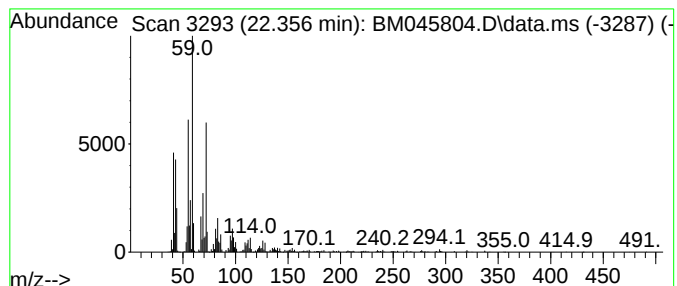
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.356	22.36 ng/ul	2402330	Chrysene-d12	21.068

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	89
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3			9-Octadecenamide	281	C18H35NO	003322-62-1	81
4			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	72
5			Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
9-Octadecenit...	20.821	2.0	ng/ul	218706	5	21.068	2148580	20.0
13-Docosenamide...	22.356	22.4	ng/ul	2402330	5	21.068	2148580	20.0