

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
 Data File : BP021465.D
 Acq On : 09 Aug 2024 09:59
 Operator : RC/JU
 Sample : P3480-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0AA3

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021465.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.146	23	27	39	rVB	31090	50220	1.25%	0.137%
2	3.599	101	104	120	rBV	56355	147812	3.69%	0.402%
3	3.870	146	150	156	rVB4	4872	8685	0.22%	0.024%
4	4.640	275	281	286	rVB10	2914	5298	0.13%	0.014%
5	6.887	655	663	671	rBV2	46450	158235	3.95%	0.431%
6	6.963	671	676	688	rVV	237003	479854	11.97%	1.306%
7	7.175	706	712	730	rBV	293586	675787	16.86%	1.839%
8	7.605	778	785	809	rBV	405839	838011	20.91%	2.280%
9	8.369	906	915	942	rBV	155864	657476	16.41%	1.789%
10	8.534	942	943	949	rVB6	3171	5927	0.15%	0.016%
11	8.787	979	986	1010	rBV	361487	872543	21.77%	2.374%
12	9.428	1086	1095	1100	rBV10	5152	17810	0.44%	0.048%
13	9.499	1100	1107	1126	rBV	219903	692950	17.29%	1.886%
14	10.075	1198	1205	1225	rBV	241530	946970	23.63%	2.577%
15	10.375	1250	1256	1273	rVB	664610	1347230	33.62%	3.666%
16	10.581	1283	1291	1309	rBV3	187339	681305	17.00%	1.854%
17	11.034	1363	1368	1369	rBV4	4212	4415	0.11%	0.012%
18	11.216	1392	1399	1412	rBV4	36173	154273	3.85%	0.420%
19	12.028	1529	1537	1542	rBV6	6689	16793	0.42%	0.046%
20	12.846	1672	1676	1681	rBV7	2538	4008	0.10%	0.011%
21	13.116	1717	1722	1735	rBV4	8122	22142	0.55%	0.060%
22	13.681	1810	1818	1838	rBV	865055	1831281	45.70%	4.983%
23	13.945	1855	1863	1883	rBV	1117587	1998563	49.87%	5.438%
24	14.245	1908	1914	1934	rBV	1121791	1894558	47.28%	5.155%
25	14.498	1952	1957	1961	rVB8	3653	6978	0.17%	0.019%
26	14.722	1989	1995	1997	rBV4	20936	40114	1.00%	0.109%
27	14.757	1997	2001	2009	rVB5	15878	43522	1.09%	0.118%
28	15.116	2058	2062	2067	rBV8	3047	5762	0.14%	0.016%
29	15.275	2083	2089	2106	rBV	1385128	2640228	65.89%	7.184%
30	15.410	2109	2112	2117	rVB6	14772	22291	0.56%	0.061%
31	15.475	2117	2123	2139	rBV2	244350	772846	19.29%	2.103%
32	15.987	2206	2210	2214	rBV6	9583	14766	0.37%	0.040%
33	16.334	2264	2269	2272	rBV7	3925	7783	0.19%	0.021%
34	16.634	2317	2320	2322	rBV4	3474	4969	0.12%	0.014%
35	16.975	2375	2378	2384	rVB8	10154	17157	0.43%	0.047%
36	17.069	2387	2394	2405	rBV	1402547	2321415	57.93%	6.317%

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37	17.181	2407	2413	2434	rVB	1857776	3091402	77.14%	8.412%
38	17.757	2507	2511	2515	rVB4	13467	18059	0.45%	0.049%
39	17.998	2547	2552	2562	rBV	151290	279910	6.98%	0.762%
40	19.110	2737	2741	2747	rBV2	26531	46794	1.17%	0.127%
41	19.198	2752	2756	2767	rVV	29990	64693	1.61%	0.176%
42	19.333	2774	2779	2786	rBV3	59190	107266	2.68%	0.292%
43	19.563	2812	2818	2839	rBV	2535308	3972852	99.14%	10.810%
44	21.157	3085	3089	3097	rBV5	92338	175802	4.39%	0.478%
45	21.522	3145	3151	3167	rBV	1549222	2653886	66.23%	7.221%
46	24.568	3660	3669	3686	rVB2	1315028	4007318	100.00%	10.904%
47	24.786	3699	3706	3730	rVB	846631	2922753	72.94%	7.953%

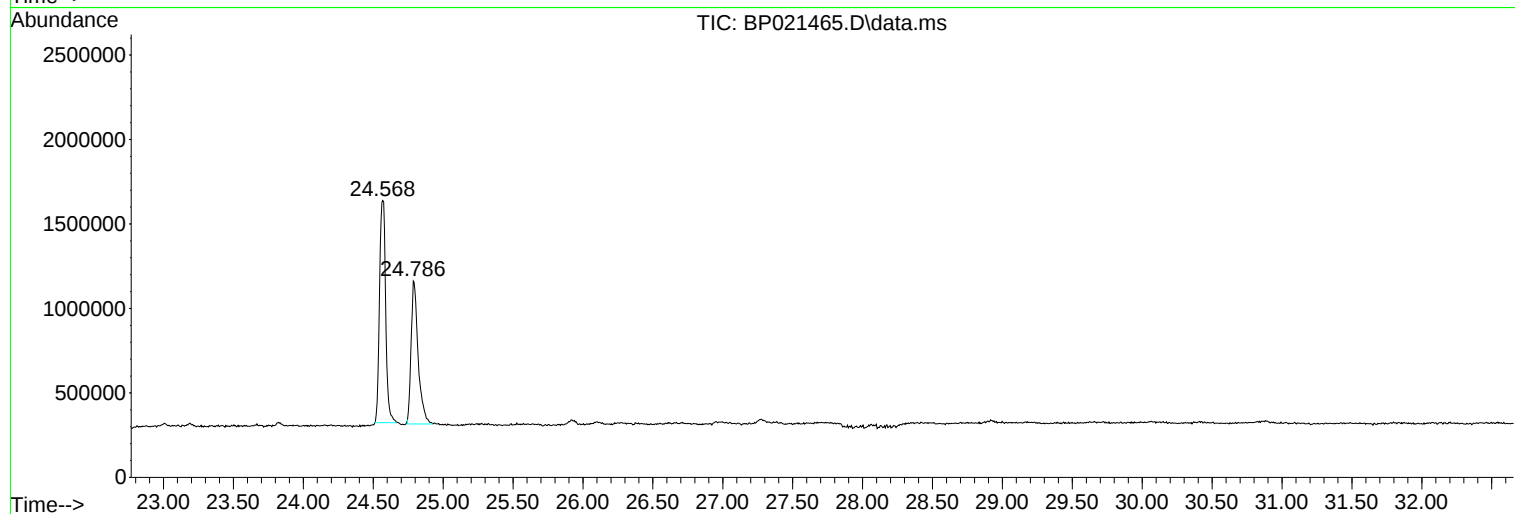
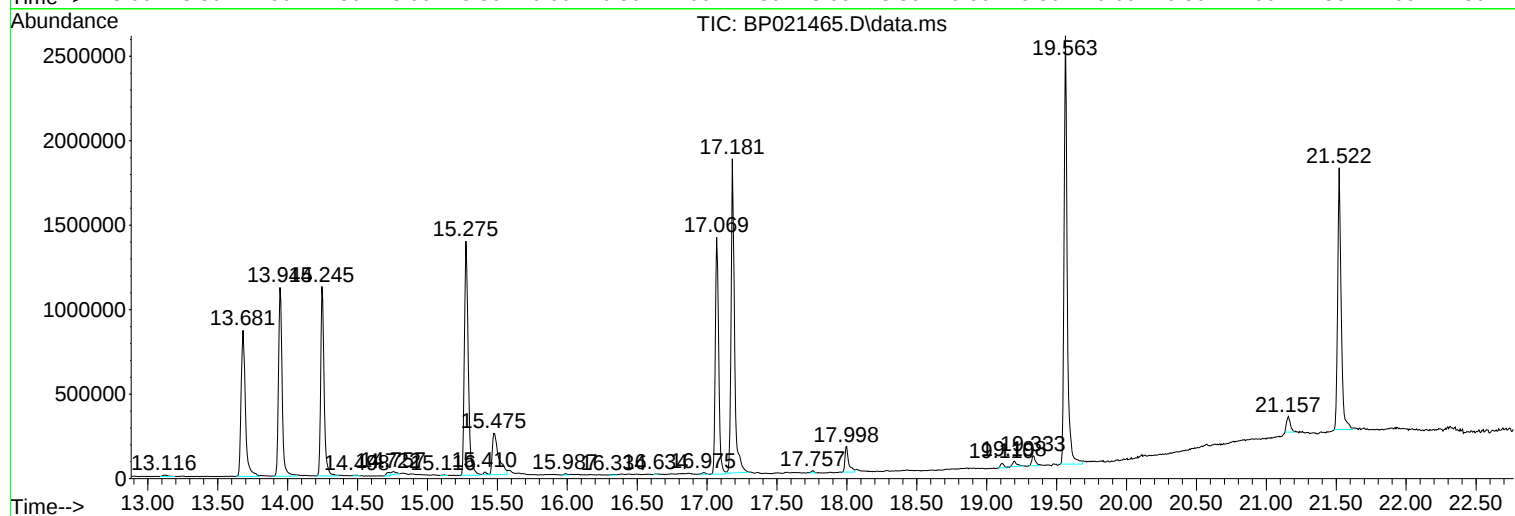
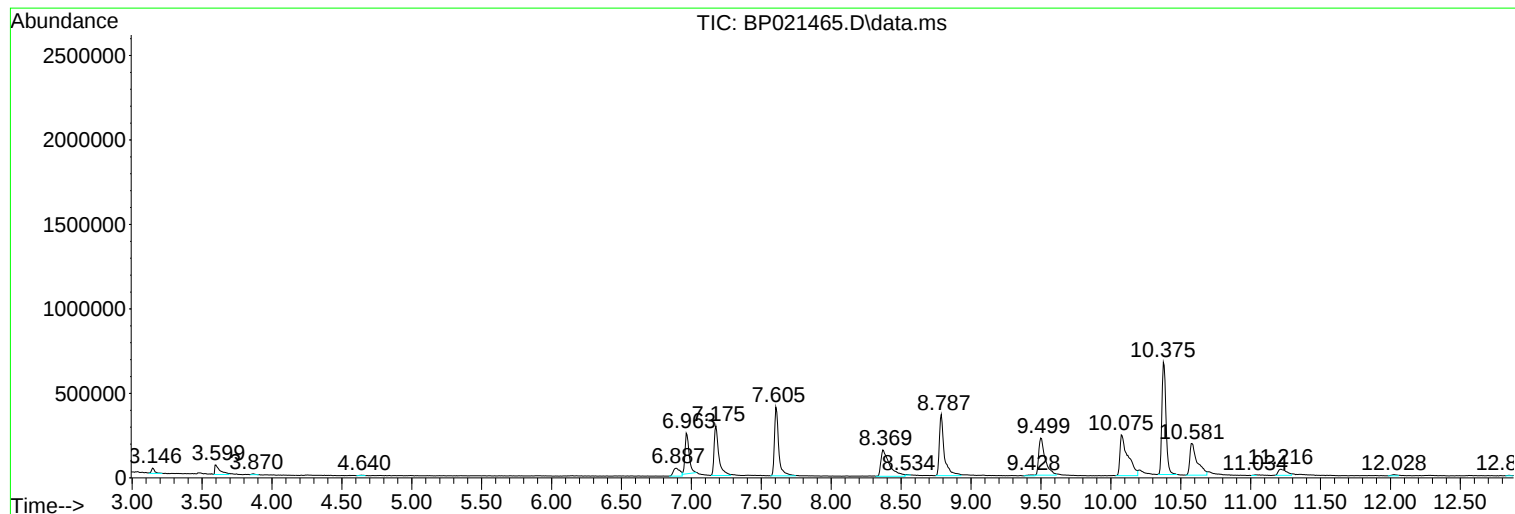
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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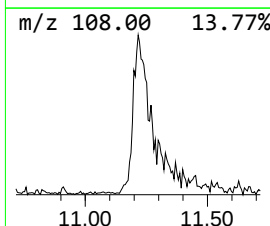
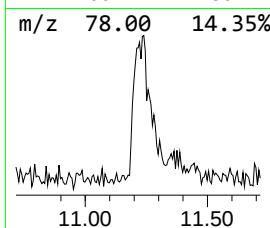
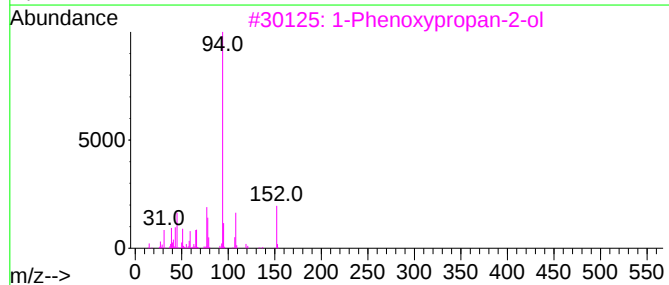
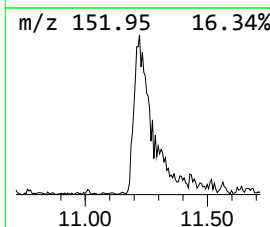
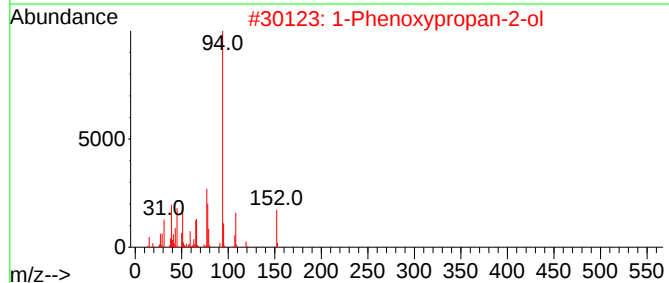
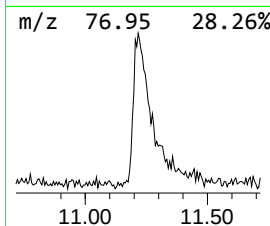
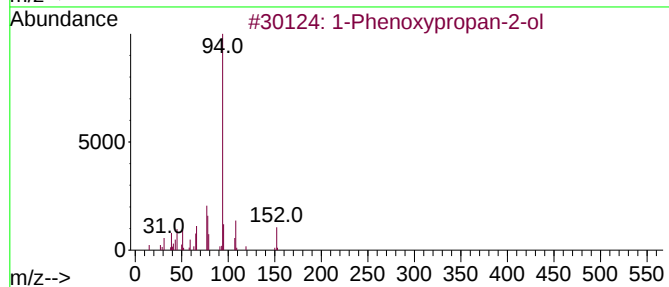
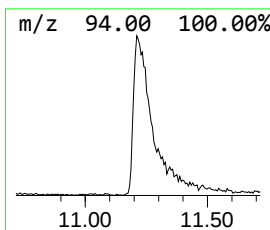
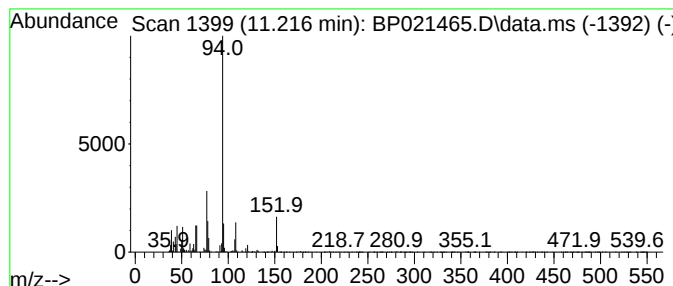
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.216	2.29 ng/ul	154273	Naphthalene-d8	10.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	86
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	83
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



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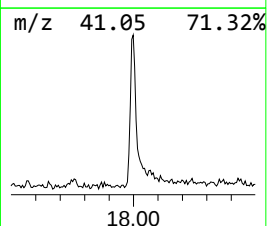
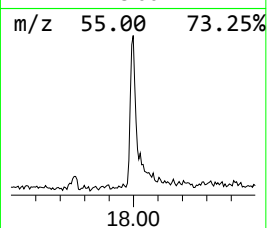
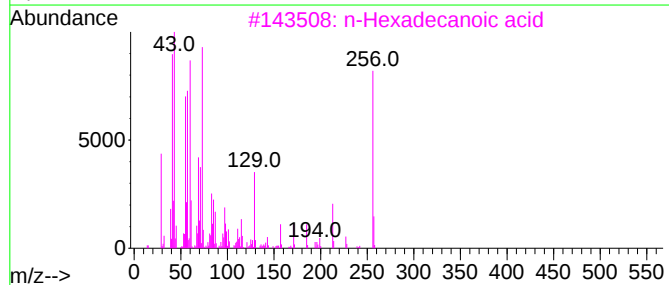
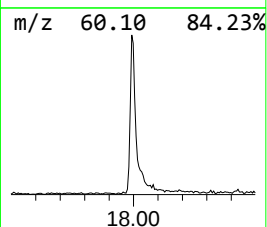
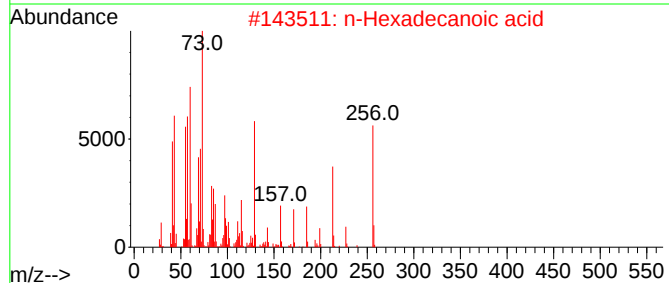
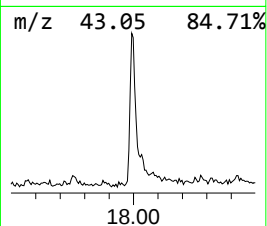
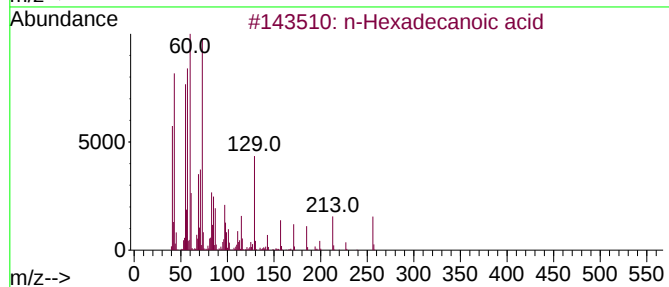
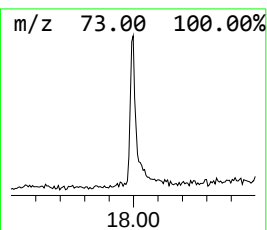
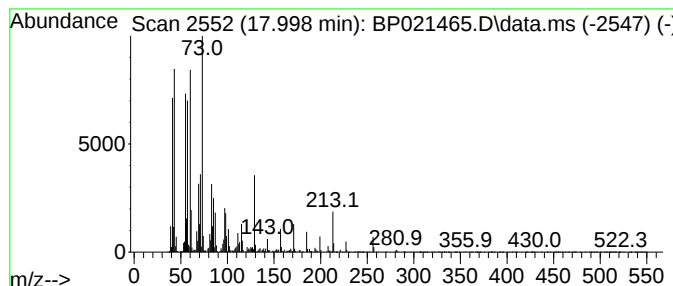
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.998	2.41 ng/ul	279910	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



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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
1-Phenoxypropan...	11.216	2.3	ng/u1	154273	2	10.375	1347230	20.0
n-Hexadecanoic ...	17.998	2.4	ng/u1	279910	4	17.069	2321420	20.0