

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
 Data File : BP021012.D
 Acq On : 17 Jul 2024 07:00
 Operator : MA/JU
 Sample : P3178-11
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4BT6

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: BP021012.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	40	45	54	rVB	36034	51709	1.45%	0.106%
2	3.528	89	92	104	rBV	72735	122065	3.43%	0.250%
3	3.681	116	118	127	rBV	44506	78420	2.20%	0.161%
4	3.764	129	132	138	rVB3	13788	17259	0.48%	0.035%
5	3.969	163	167	173	rVB	22551	31240	0.88%	0.064%
6	4.381	235	237	243	rVB	9750	13988	0.39%	0.029%
7	4.469	249	252	257	rBV3	5528	8363	0.23%	0.017%
8	4.522	257	261	265	rVB2	11264	18191	0.51%	0.037%
9	4.746	295	299	306	rBV10	5923	11179	0.31%	0.023%
10	4.863	314	319	329	rVB	54924	80045	2.25%	0.164%
11	5.828	475	483	493	rBV5	8637	28414	0.80%	0.058%
12	6.805	638	649	655	rVB10	7705	25960	0.73%	0.053%
13	6.899	655	665	681	rBV	604821	1108488	31.11%	2.270%
14	7.046	683	690	704	rVV	517621	824529	23.14%	1.689%
15	7.246	716	724	732	rBV	651907	1126483	31.61%	2.307%
16	7.693	788	800	810	rBV	830552	1300000	36.48%	2.662%
17	7.775	812	814	822	rVB6	11159	17175	0.48%	0.035%
18	8.187	879	884	891	rBV6	5233	9018	0.25%	0.018%
19	8.310	901	905	911	rVB9	4429	7028	0.20%	0.014%
20	8.410	911	922	936	rBV	725551	1351415	37.92%	2.768%
21	8.522	937	941	951	rVB3	14143	33678	0.95%	0.069%
22	8.804	984	989	991	rBV	29033	46362	1.30%	0.095%
23	8.851	991	997	1012	rVB	609250	1066663	29.93%	2.184%
24	9.199	1051	1056	1063	rBV3	10709	19876	0.56%	0.041%
25	9.393	1083	1089	1102	rVB	59071	102141	2.87%	0.209%
26	9.551	1107	1116	1129	rBV	494854	920338	25.83%	1.885%
27	9.922	1173	1179	1185	rVB5	5969	13924	0.39%	0.029%
28	10.098	1201	1209	1228	rBV	750037	1449510	40.68%	2.968%
29	10.457	1262	1270	1276	rBV	1136192	1792802	50.31%	3.672%
30	10.604	1288	1295	1319	rBV	476412	1040038	29.19%	2.130%
31	10.993	1354	1361	1370	rBV9	16228	43485	1.22%	0.089%
32	11.204	1389	1397	1408	rBV2	100430	241664	6.78%	0.495%
33	11.669	1471	1476	1482	rVB9	3642	7912	0.22%	0.016%
34	11.857	1503	1508	1514	rBV9	4788	10287	0.29%	0.021%
35	12.034	1531	1538	1548	rVB3	12396	24550	0.69%	0.050%
36	12.587	1629	1632	1639	rBV9	3233	7615	0.21%	0.016%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
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37	13.104	1715	1720	1727	rBV6	6702	10166	0.29%	0.021%
38	13.263	1742	1747	1748	rBV5	5421	7643	0.21%	0.016%
39	13.498	1777	1787	1789	rBV10	4530	11859	0.33%	0.024%
40	13.610	1801	1806	1812	rVB10	8897	18197	0.51%	0.037%

41	13.722	1814	1825	1837	rBV	1229712	2086289	58.55%	4.273%
42	13.816	1837	1841	1848	rVV8	17725	28807	0.81%	0.059%
43	13.881	1848	1852	1859	rVB10	5630	11380	0.32%	0.023%
44	13.992	1861	1871	1889	rBV	1667069	2610100	73.25%	5.345%
45	14.134	1890	1895	1899	rVV4	24881	37597	1.06%	0.077%

46	14.181	1900	1903	1908	rVB6	6280	10202	0.29%	0.021%
47	14.304	1916	1924	1931	rBV	1752028	2481520	69.64%	5.082%
48	14.369	1931	1935	1941	rVV3	56895	95933	2.69%	0.196%
49	14.422	1941	1944	1949	rVB6	11409	14153	0.40%	0.029%
50	14.534	1955	1963	1965	rBV4	90461	204049	5.73%	0.418%

51	14.575	1966	1970	1987	rVB	324082	806361	22.63%	1.651%
52	14.763	1999	2002	2011	rVB2	52947	79090	2.22%	0.162%
53	15.039	2046	2049	2054	rVB6	7443	10802	0.30%	0.022%
54	15.110	2059	2061	2065	rVB2	10455	10740	0.30%	0.022%
55	15.163	2065	2070	2080	rVB2	7779	19837	0.56%	0.041%

56	15.257	2082	2086	2089	rBV5	7134	12315	0.35%	0.025%
57	15.316	2090	2096	2105	rBV	1912701	3192343	89.59%	6.538%
58	15.475	2116	2123	2132	rBV	469572	807353	22.66%	1.653%
59	15.904	2191	2196	2202	rBV8	23975	46579	1.31%	0.095%
60	15.963	2204	2206	2209	rBV4	5172	7675	0.22%	0.016%

61	16.063	2216	2223	2238	rVB4	32482	94627	2.66%	0.194%
62	16.639	2317	2321	2325	rVB6	14645	23092	0.65%	0.047%
63	16.857	2356	2358	2360	rBV2	9617	9288	0.26%	0.019%
64	16.986	2376	2380	2383	rBV	36987	47062	1.32%	0.096%
65	17.022	2383	2386	2394	rVB6	27273	42791	1.20%	0.088%

66	17.122	2397	2403	2407	rBV	2315383	3021449	84.79%	6.188%
67	17.157	2407	2409	2413	rVV	207239	243560	6.83%	0.499%
68	17.216	2413	2419	2429	rVV	2197932	3368543	94.53%	6.899%
69	17.292	2429	2432	2435	rVB2	34781	42094	1.18%	0.086%
70	17.575	2478	2480	2484	rVB	30002	32243	0.90%	0.066%

71	17.739	2501	2508	2513	rVB	137608	251822	7.07%	0.516%
72	17.851	2525	2527	2533	rVB6	25000	37015	1.04%	0.076%
73	18.045	2554	2560	2569	rBV2	256680	521526	14.64%	1.068%
74	18.222	2586	2590	2595	rBV2	154264	231762	6.50%	0.475%
75	18.280	2597	2600	2604	rVV2	89750	133443	3.74%	0.273%

76	18.327	2605	2608	2614	rVB4	65166	98163	2.75%	0.201%
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77	18.516	2636	2640	2644	rBV2	73377	107085	3.01%	0.219%
78	18.698	2668	2671	2675	rVB2	68110	77729	2.18%	0.159%
79	19.010	2721	2724	2725	rBV	36636	41364	1.16%	0.085%
80	19.080	2733	2736	2738	rBV	53129	63366	1.78%	0.130%
81	19.251	2761	2765	2771	rVB	365347	532884	14.95%	1.091%
82	19.380	2783	2787	2796	rBV4	93221	223745	6.28%	0.458%
83	19.598	2818	2824	2839	rBV	2221455	3563477	100.00%	7.298%
84	21.210	3093	3098	3108	rVB2	487093	942023	26.44%	1.929%
85	21.551	3149	3156	3162	rBV2	1840775	3501209	98.25%	7.170%
86	23.933	3557	3561	3569	rVB	294893	585643	16.43%	1.199%
87	24.615	3671	3677	3685	rVB2	853212	1918228	53.83%	3.928%
88	24.845	3707	3716	3724	rBV	1413767	3402020	95.47%	6.967%

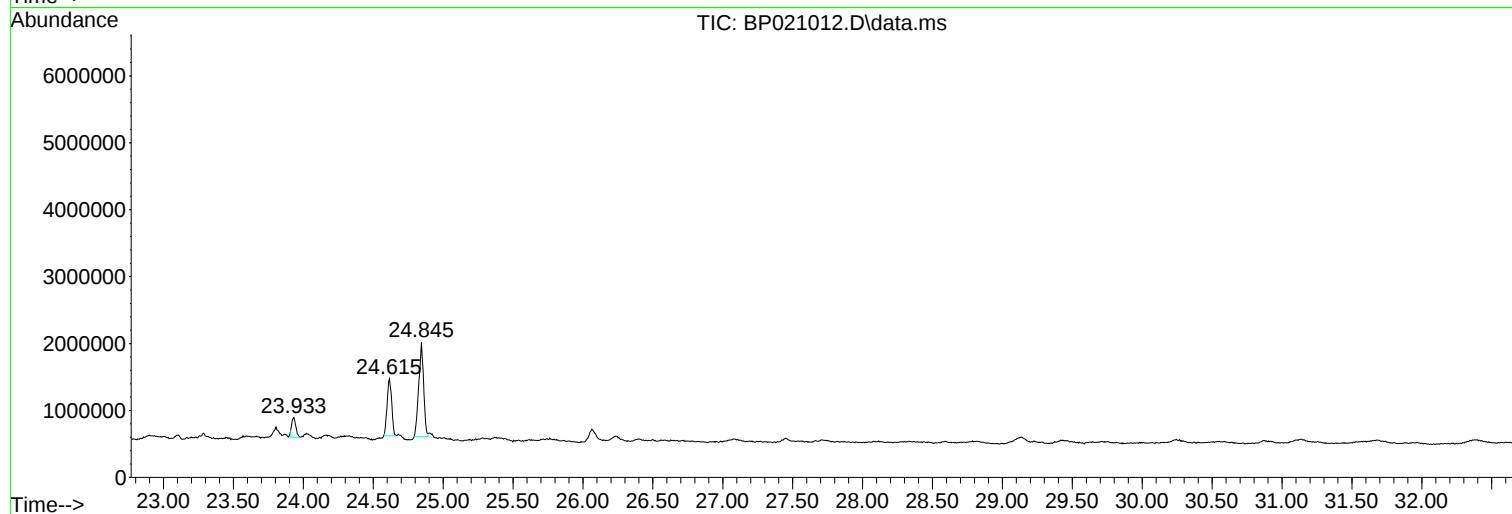
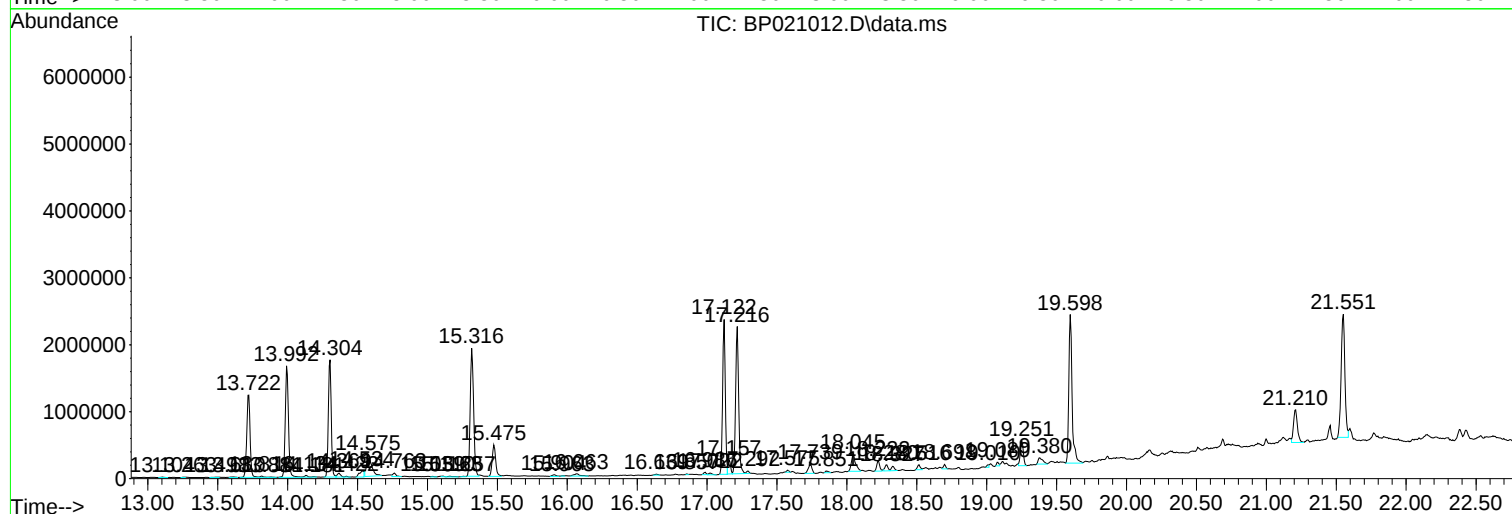
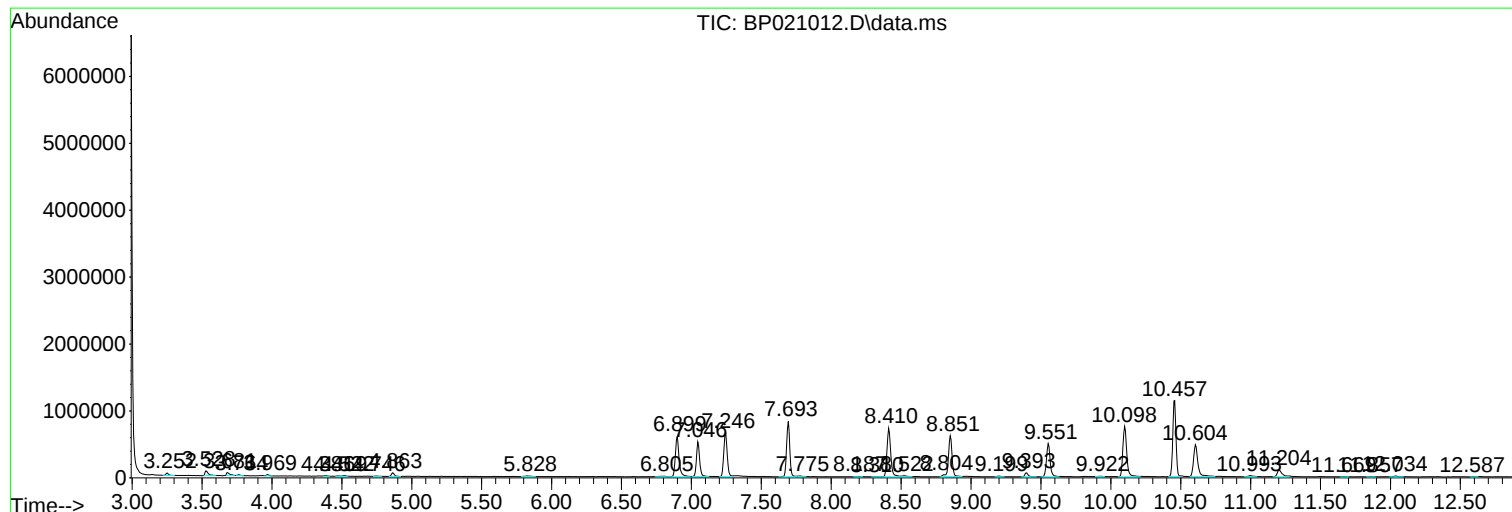
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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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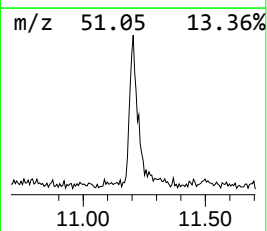
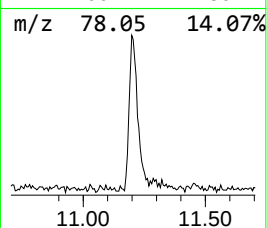
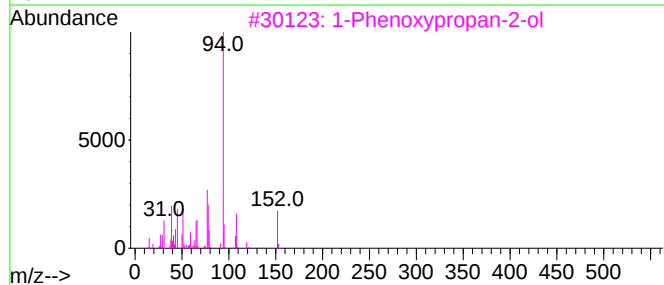
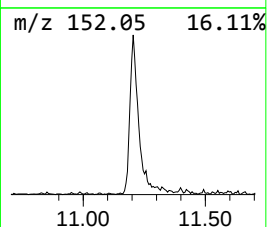
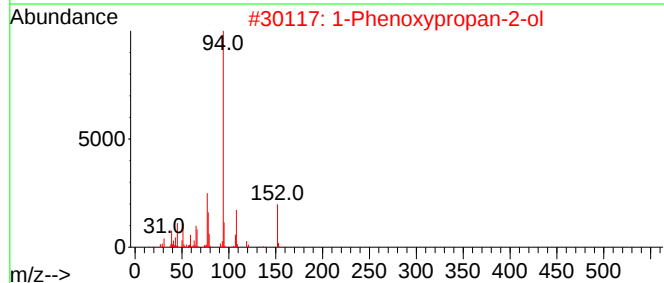
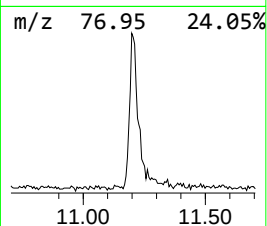
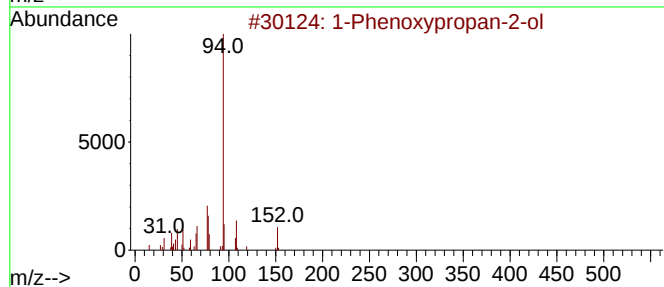
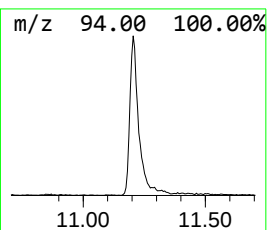
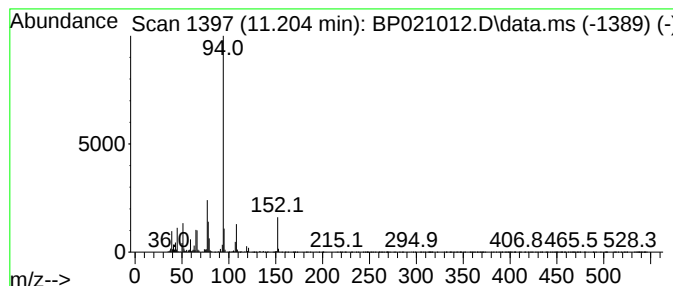
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.204	2.70 ng/ul	241664	Naphthalene-d8	10.451

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
5			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64



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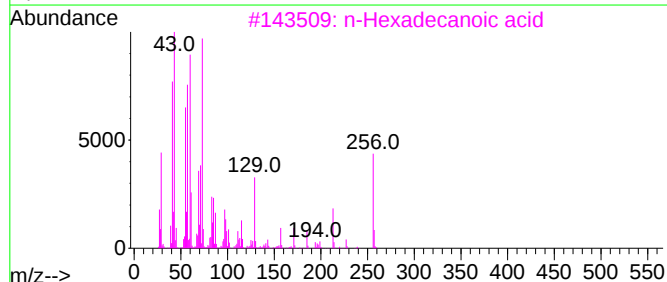
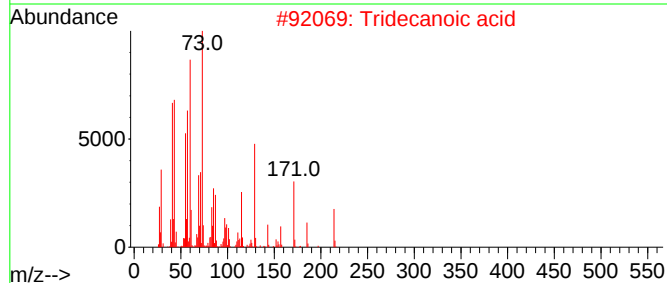
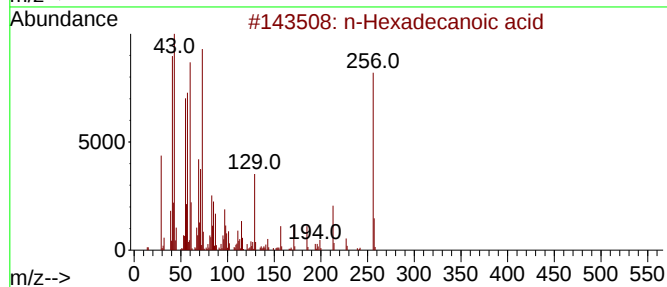
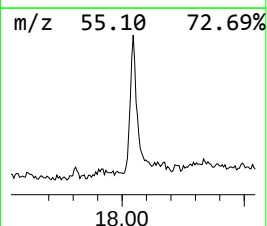
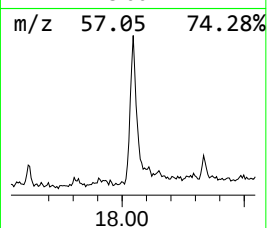
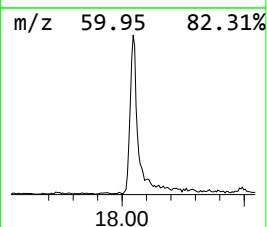
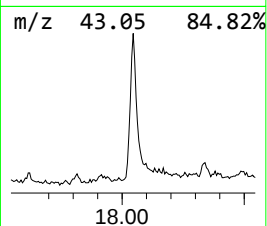
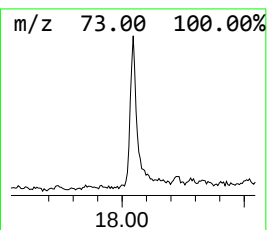
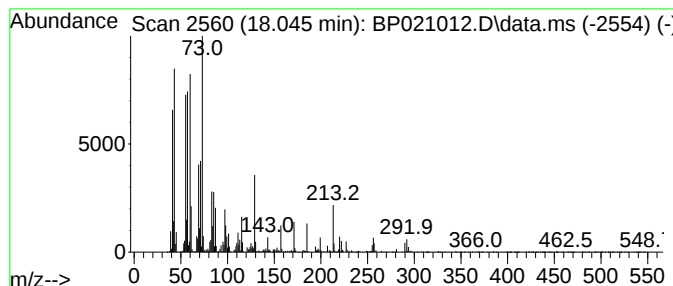
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	3.45 ng/ul	521526	Phenanthrene-d10	17.122

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	95
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



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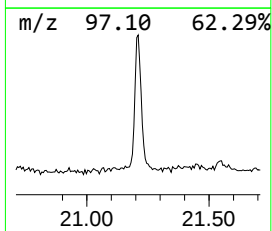
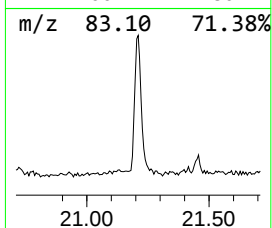
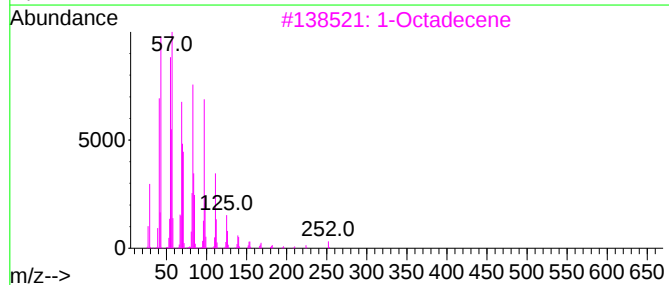
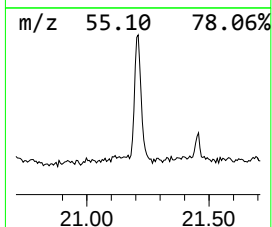
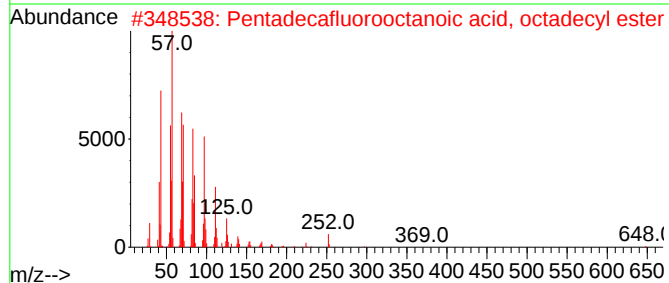
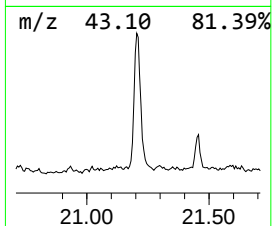
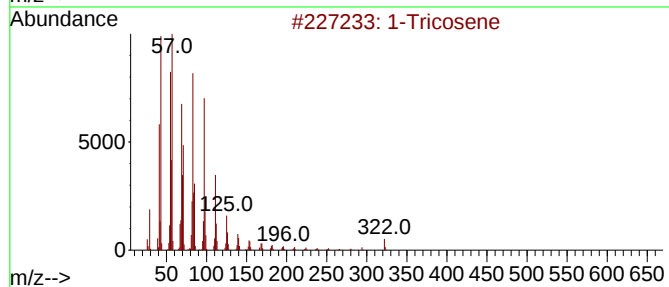
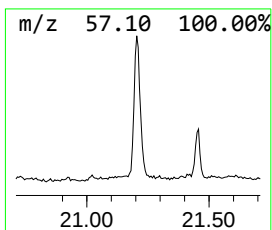
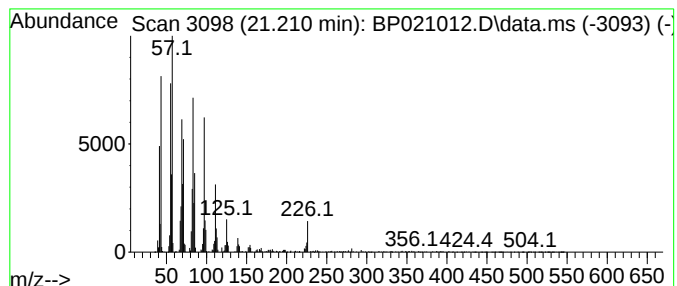
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.210	5.38 ng/ul	942023	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene			322	C23H46	018835-32-0	98
2	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	95
3	1-Octadecene			252	C18H36	000112-88-9	94
4	Cyclohexadecane			224	C16H32	000295-65-8	94
5	2- Chloropropionic acid, hexadec...			332	C19H37ClO2	086711-81-1	93



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ClientSampleId :
A4BT6

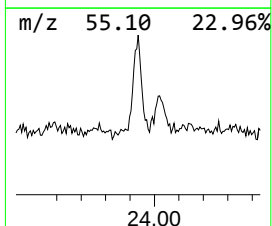
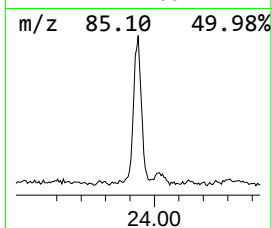
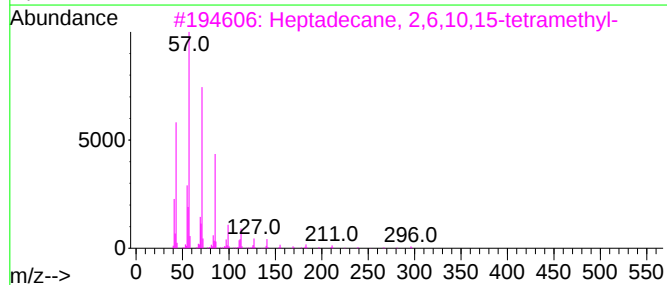
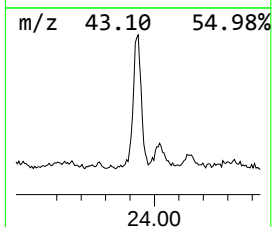
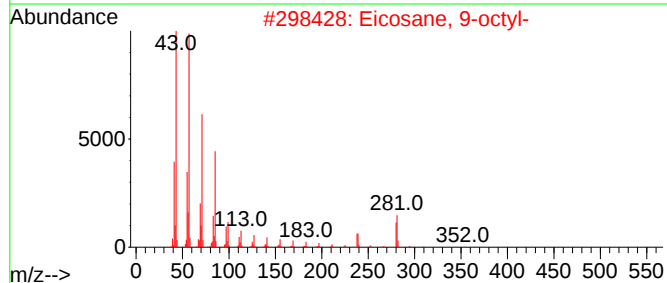
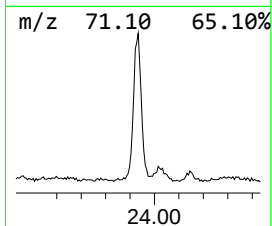
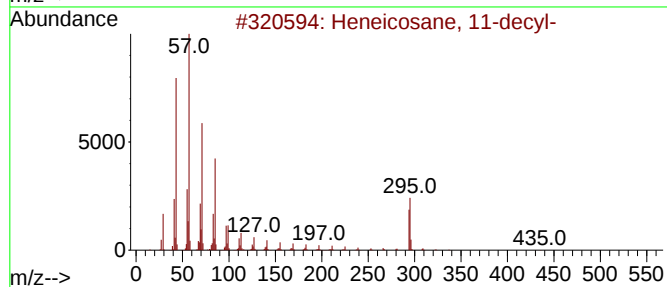
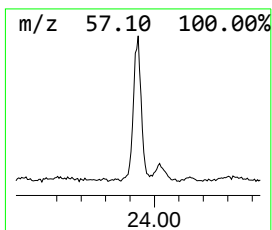
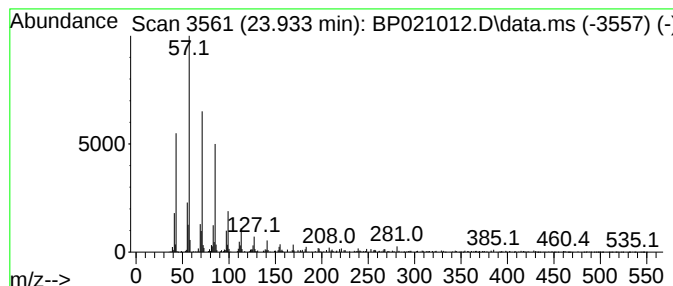
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.933	3.44 ng/ul	585643	Perylene-d12	24.845

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	93
2			Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
3			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
4			Octacosane	394	C28H58	000630-02-4	90
5			Pentadecane, 7-(bromomethyl)-	304	C16H33Br	052997-43-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071524\
Data File : BP021012.D
Acq On : 17 Jul 2024 07:00
Operator : MA/JU
Sample : P3178-11
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BT6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
1-Phenoxypropan...	11.204	2.7	ng/u1	241664	2	10.451	1792800	20.0
n-Hexadecanoic ...	18.045	3.5	ng/u1	521526	4	17.122	3021450	20.0
(DEL) Alkane: S...	21.210	5.4	ng/u1	942023	5	21.551	3501210	20.0
(DEL) Alkane: S...	23.933	3.4	ng/u1	585643	6	24.845	3402020	20.0