

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010020.D
 Acq On : 21 Apr 2022 18:26
 Operator : CG/JU
 Sample : N2373-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFD2

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-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010020.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.799	118	121	132	rBV	40051	71791	1.65%	0.181%
2	3.893	134	137	141	rVB3	21167	27323	0.63%	0.069%
3	7.098	675	682	693	rVB	246434	410433	9.41%	1.032%
4	7.263	704	710	718	rBV	190334	321462	7.37%	0.808%
5	7.469	735	745	759	rBV	241468	413776	9.49%	1.040%
6	7.939	818	825	835	rBV	348564	584037	13.39%	1.468%
7	8.251	872	878	889	rVB	194175	320238	7.34%	0.805%
8	8.645	937	945	957	rBV	318532	539849	12.38%	1.357%
9	8.763	959	965	970	rVV	21683	37452	0.86%	0.094%
10	8.810	970	973	979	rVB6	11335	18618	0.43%	0.047%
11	9.098	1012	1022	1030	rBV	261769	470547	10.79%	1.183%
12	9.551	1094	1099	1105	rVB4	24372	40900	0.94%	0.103%
13	9.822	1138	1145	1153	rBV	215153	379355	8.70%	0.954%
14	10.010	1169	1177	1184	rBV7	14134	28050	0.64%	0.071%
15	10.363	1230	1237	1247	rBV	359197	633843	14.53%	1.594%
16	10.522	1256	1264	1274	rBV7	9882	25935	0.59%	0.065%
17	10.745	1294	1302	1309	rBV	524170	923029	21.16%	2.321%
18	10.881	1316	1325	1338	rBV	184813	375494	8.61%	0.944%
19	11.292	1389	1395	1403	rVB8	11351	21460	0.49%	0.054%
20	11.875	1489	1494	1501	rBV6	10481	18421	0.42%	0.046%
21	12.180	1540	1546	1551	rBV2	10833	18751	0.43%	0.047%
22	13.410	1750	1755	1761	rBV2	19230	25325	0.58%	0.064%
23	13.757	1808	1814	1819	rVB8	8523	18070	0.41%	0.045%
24	13.892	1832	1837	1844	rBV4	19276	39878	0.91%	0.100%
25	13.974	1844	1851	1858	rBV	833516	1210899	27.76%	3.045%
26	14.263	1889	1900	1915	rBV	835489	1313625	30.12%	3.303%
27	14.569	1941	1952	1961	rBV	922390	1392872	31.93%	3.502%
28	14.757	1978	1984	1993	rBV	256569	444567	10.19%	1.118%
29	14.827	1994	1996	2000	rVV4	19882	27229	0.62%	0.068%
30	14.892	2004	2007	2015	rVV3	17259	43501	1.00%	0.109%
31	14.974	2015	2021	2028	rVV6	17577	40201	0.92%	0.101%
32	15.045	2029	2033	2038	rVV3	16980	24129	0.55%	0.061%
33	15.186	2050	2057	2063	rBV7	13700	33828	0.78%	0.085%
34	15.363	2080	2087	2088	rBV7	9777	17625	0.40%	0.044%
35	15.563	2114	2121	2127	rVV	1181580	1710232	39.21%	4.300%
36	15.616	2127	2130	2134	rVB4	24299	35020	0.80%	0.088%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

37	15.674	2134	2140	2150	rBV	368252	543020	12.45%	1.365%
38	15.804	2158	2162	2170	rVB6	15577	36104	0.83%	0.091%
39	15.939	2174	2185	2192	rBV6	14465	45011	1.03%	0.113%
40	16.016	2194	2198	2203	rVV8	7888	18951	0.43%	0.048%
41	16.063	2203	2206	2212	rVB7	9722	17838	0.41%	0.045%
42	16.139	2214	2219	2228	rVB10	17708	37426	0.86%	0.094%
43	16.298	2240	2246	2251	rVV2	54709	83028	1.90%	0.209%
44	16.627	2295	2302	2307	rBV4	14169	28000	0.64%	0.070%
45	16.674	2307	2310	2315	rVB	13946	19532	0.45%	0.049%
46	16.774	2323	2327	2333	rVB7	12784	20140	0.46%	0.051%
47	16.898	2344	2348	2352	rVV6	15278	26574	0.61%	0.067%
48	16.968	2354	2360	2371	rVB3	27244	69845	1.60%	0.176%
49	17.145	2386	2390	2393	rBV	21723	27234	0.62%	0.068%
50	17.321	2406	2420	2425	rBV	1201957	1852201	42.46%	4.657%
51	17.363	2425	2427	2432	rVV	289986	378075	8.67%	0.951%
52	17.421	2432	2437	2448	rVB2	1313025	1975185	45.28%	4.966%
53	17.521	2448	2454	2461	rBV3	141696	248503	5.70%	0.625%
54	17.604	2464	2468	2471	rVV6	23356	43285	0.99%	0.109%
55	17.639	2471	2474	2477	rVB4	17924	21430	0.49%	0.054%
56	17.904	2515	2519	2525	rVB2	37202	55517	1.27%	0.140%
57	18.074	2539	2548	2557	rVV3	210603	396371	9.09%	0.997%
58	18.192	2563	2568	2572	rBV2	227547	373581	8.56%	0.939%
59	18.233	2572	2575	2580	rVB	131628	180816	4.15%	0.455%
60	18.333	2589	2592	2594	rBV2	37448	42261	0.97%	0.106%
61	18.368	2594	2598	2602	rVV2	111262	179028	4.10%	0.450%
62	18.410	2602	2605	2614	rVB	103938	156707	3.59%	0.394%
63	18.527	2622	2625	2628	rVB5	17835	20611	0.47%	0.052%
64	18.568	2628	2632	2636	rVB4	18068	25099	0.58%	0.063%
65	18.668	2644	2649	2652	rBV2	58331	81939	1.88%	0.206%
66	18.704	2652	2655	2662	rVB3	34750	56286	1.29%	0.142%
67	18.880	2681	2685	2687	rBV4	20949	29688	0.68%	0.075%
68	18.909	2687	2690	2694	rVV	36918	54652	1.25%	0.137%
69	18.957	2694	2698	2702	rVV	46413	65898	1.51%	0.166%
70	18.992	2702	2704	2708	rVB	22171	30436	0.70%	0.077%
71	19.074	2713	2718	2722	rBV	131062	214950	4.93%	0.540%
72	19.115	2722	2725	2731	rVB4	43371	83162	1.91%	0.209%
73	19.162	2731	2733	2737	rVB	34914	37001	0.85%	0.093%
74	19.209	2737	2741	2743	rBV	50221	81051	1.86%	0.204%
75	19.327	2756	2761	2767	rVB	367614	472174	10.83%	1.187%
76	19.392	2768	2772	2774	rVV3	24238	32063	0.74%	0.081%

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77	19.427	2775	2778	2782	rVB3	30737	41334	0.95%	0.104%
78	19.515	2790	2793	2797	rVB5	22635	25505	0.58%	0.064%
79	19.568	2799	2802	2804	rBV2	17508	20838	0.48%	0.052%
80	19.651	2811	2816	2820	rBV	2011501	2833929	64.97%	7.125%
81	19.757	2830	2834	2840	rVB5	36912	60312	1.38%	0.152%
82	19.904	2857	2859	2868	rVB8	25598	39629	0.91%	0.100%
83	19.992	2870	2874	2876	rBV3	58489	85980	1.97%	0.216%
84	20.156	2893	2902	2911	rBV3	90515	254863	5.84%	0.641%
85	20.256	2916	2919	2923	rVV2	36228	47472	1.09%	0.119%
86	20.315	2925	2929	2934	rVB	86746	122359	2.81%	0.308%
87	20.451	2949	2952	2956	rBV	78946	94526	2.17%	0.238%
88	20.498	2956	2960	2964	rVB	75517	89336	2.05%	0.225%
89	20.892	3024	3027	3033	rVB2	58241	91445	2.10%	0.230%
90	21.109	3059	3064	3072	rBV3	468045	682856	15.66%	1.717%
91	21.403	3107	3114	3118	rBV	1504731	2209516	50.66%	5.555%
92	21.439	3118	3120	3131	rVB	208163	312641	7.17%	0.786%
93	22.045	3217	3223	3231	rVB4	382821	749443	17.18%	1.884%
94	23.109	3398	3404	3410	rBV	2532862	4361859	100.00%	10.967%
95	23.697	3498	3504	3510	rBV2	1084909	2046563	46.92%	5.146%
96	23.856	3524	3531	3539	rBV2	832654	1739454	39.88%	4.373%
97	24.509	3634	3642	3650	rVB	891623	1964201	45.03%	4.938%
98	24.597	3652	3657	3667	rVB3	131407	294068	6.74%	0.739%
99	26.403	3957	3964	3980	rVB3	301809	1003136	23.00%	2.522%
100	26.986	4058	4063	4074	rVB9	129986	383554	8.79%	0.964%

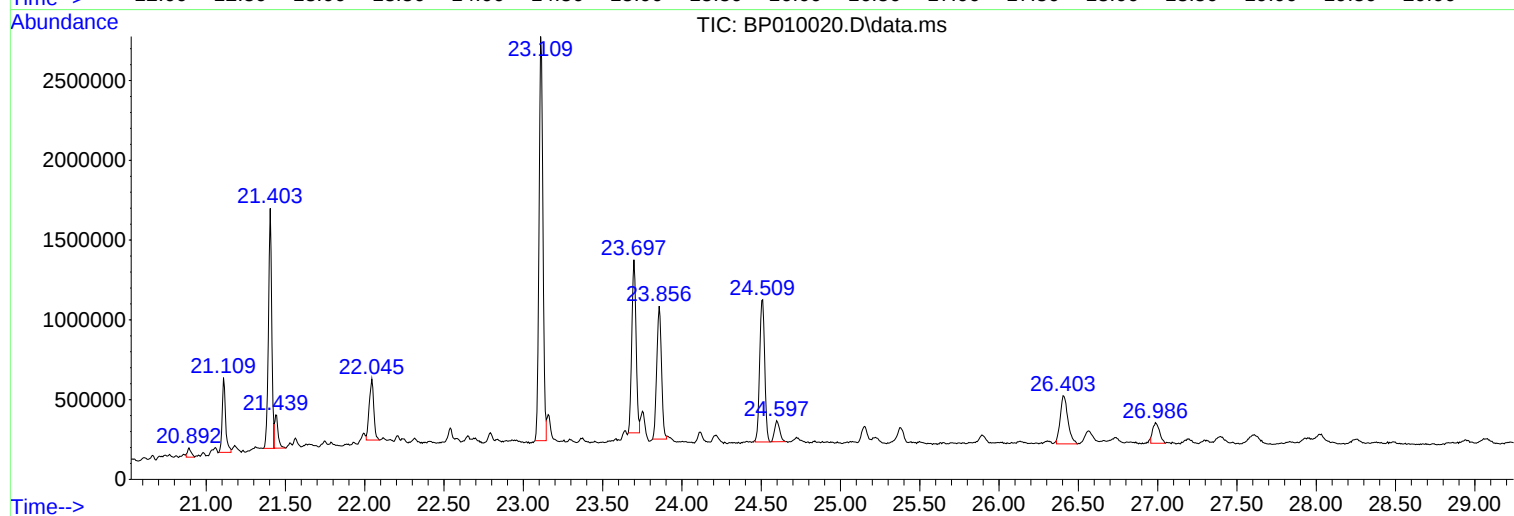
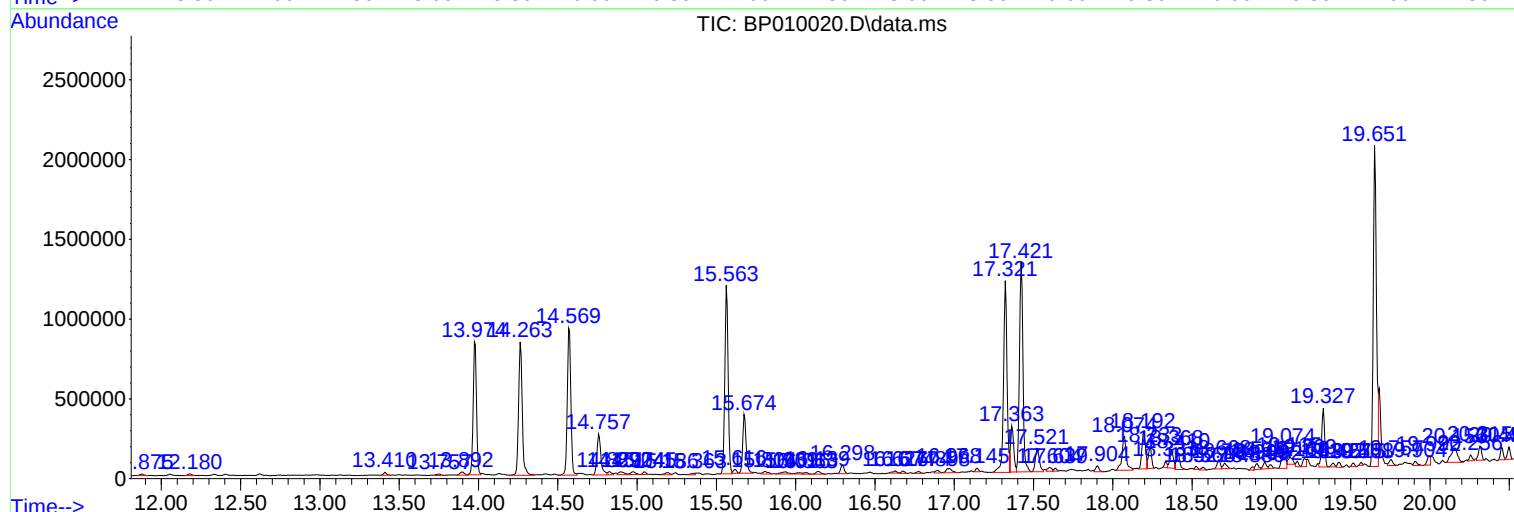
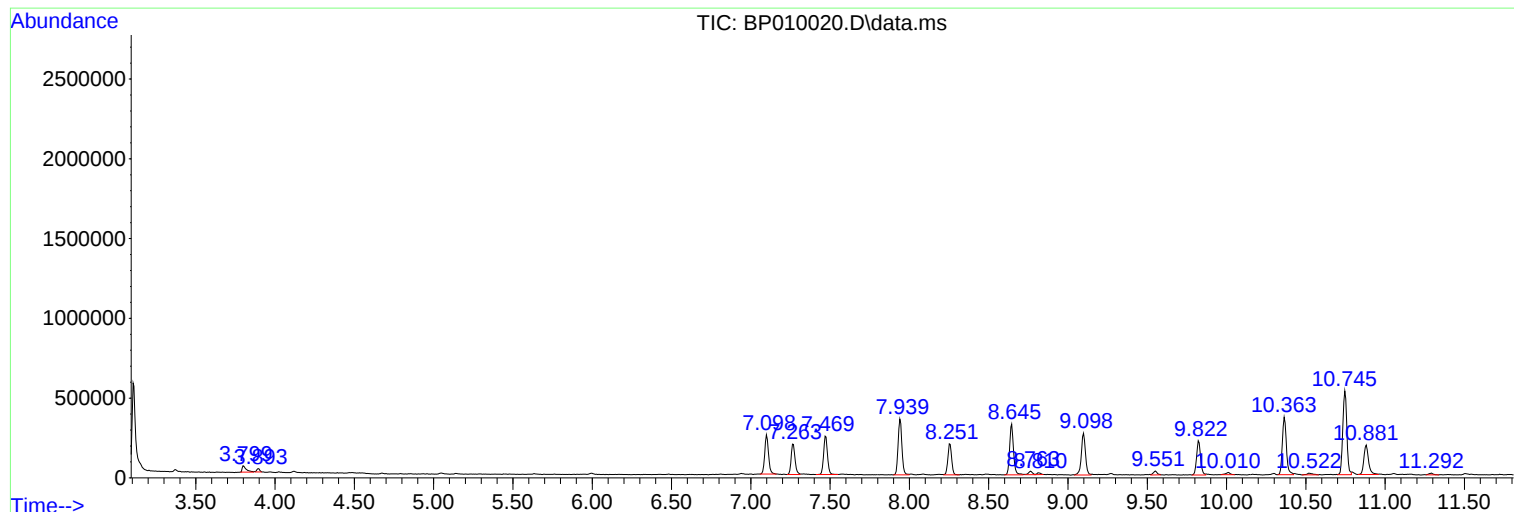
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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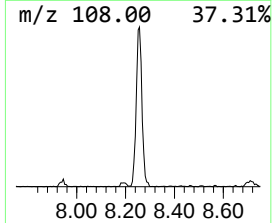
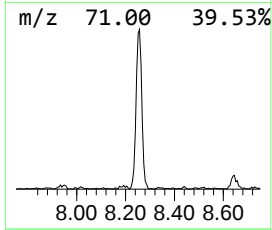
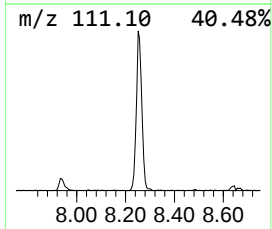
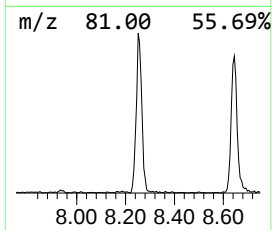
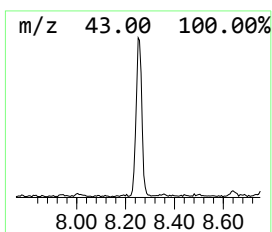
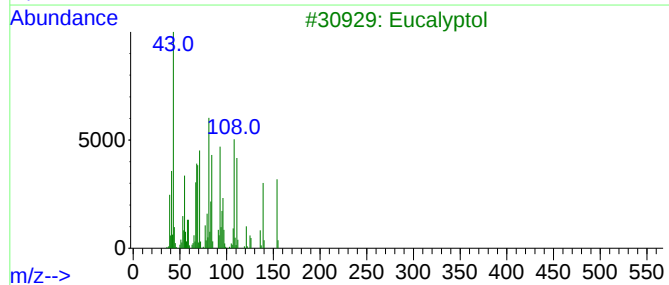
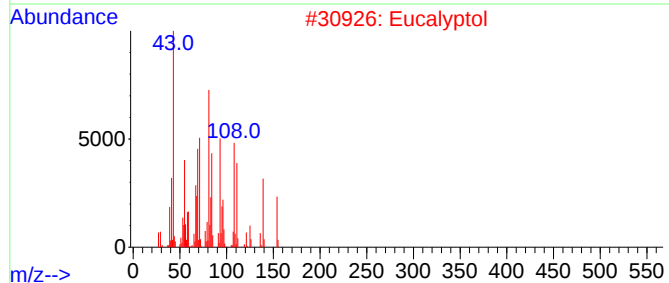
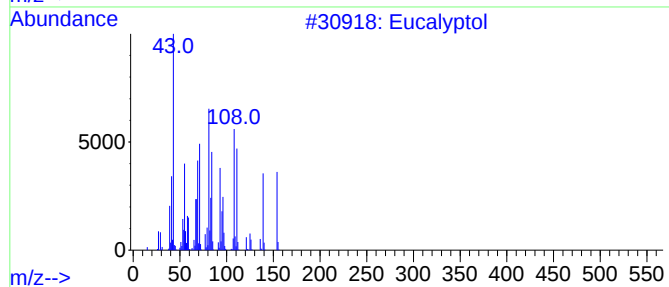
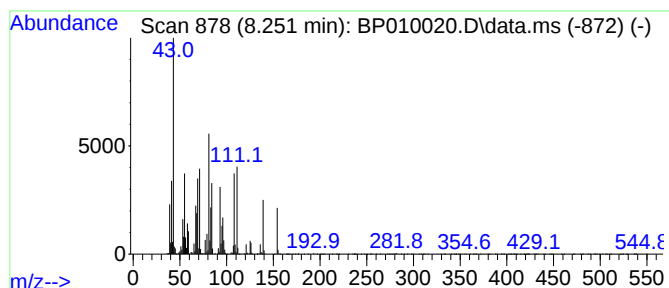
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Eucalyptol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.251	10.97 ng/ul	320238	1,4-Dichlorobenzene-d4	7.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol			154	C10H18O	000470-82-6	98
2	Eucalyptol			154	C10H18O	000470-82-6	96
3	Eucalyptol			154	C10H18O	000470-82-6	91
4	Eucalyptol			154	C10H18O	000470-82-6	76
5	Eucalyptol			154	C10H18O	000470-82-6	50



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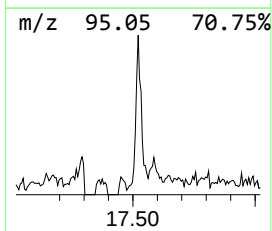
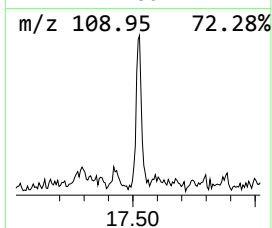
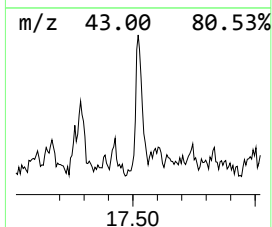
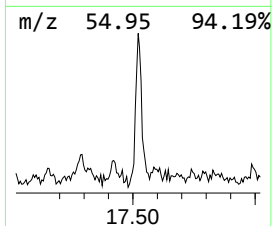
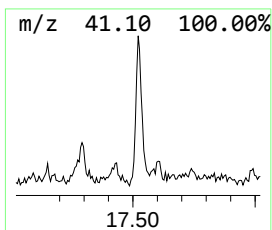
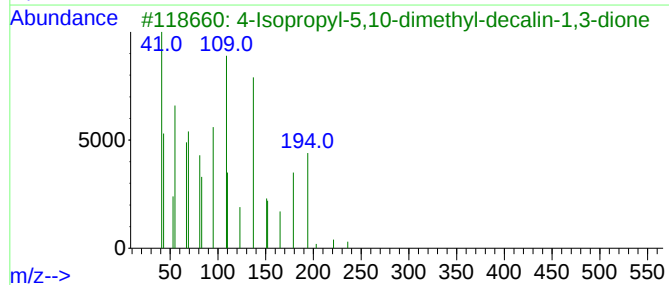
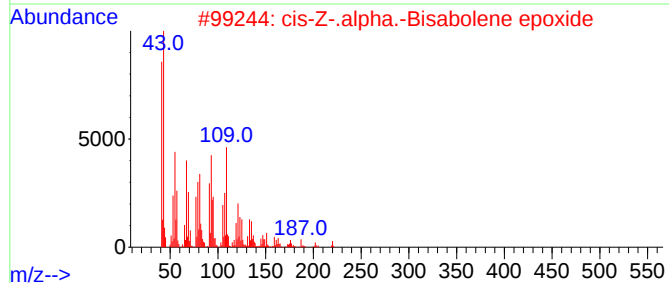
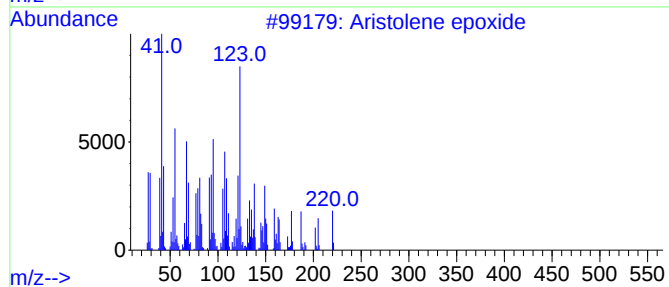
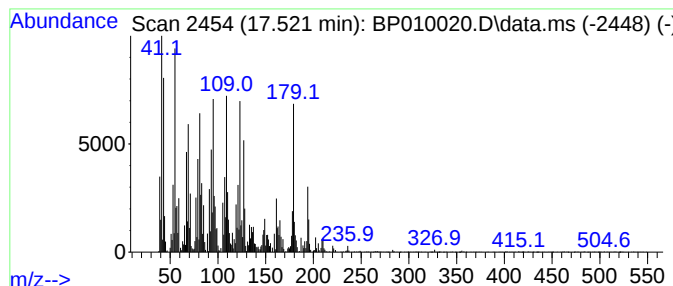
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.521	2.68 ng/ul	248503	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Aristolene epoxide	220	C15H24O	1000151-48-9	46
2			cis-Z-.alpha.-Bisabolene epoxide	220	C15H24O	1000131-71-2	46
3			4-Isopropyl-5,10-dimethyl-decali...	236	C15H24O2	291278-94-9	45
4			2(1H)-Naphthalenone, octahydro-4...	180	C12H20O	051557-64-3	35
5			2-(1-Hydroxycycloheptyl)-furan	180	C11H16O2	115754-89-7	25



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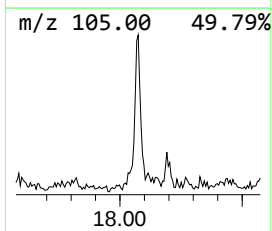
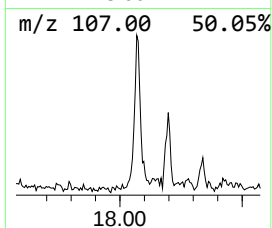
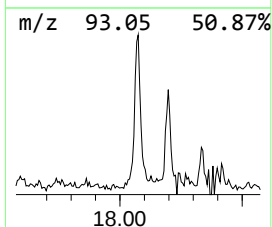
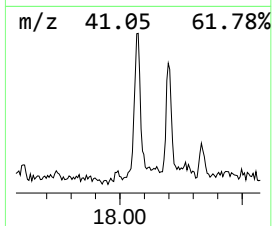
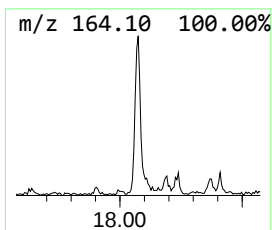
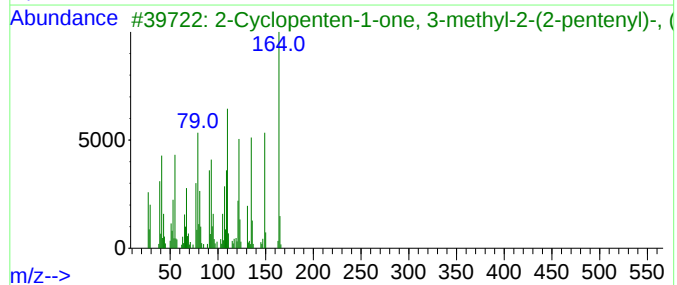
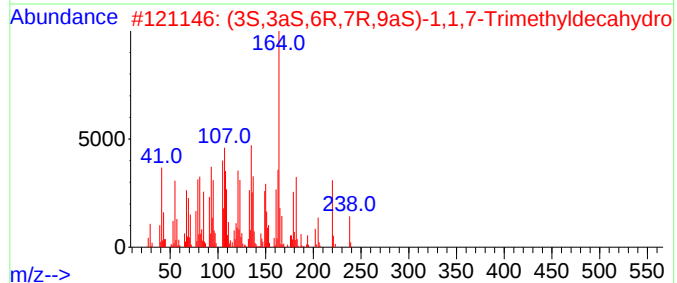
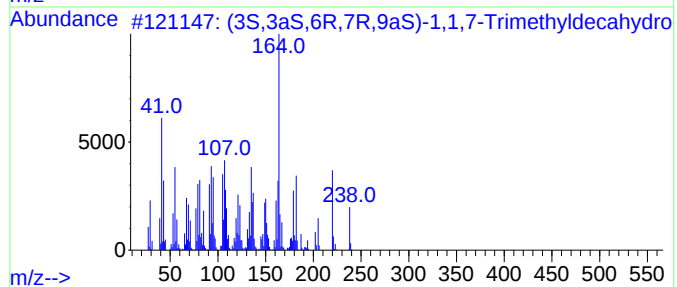
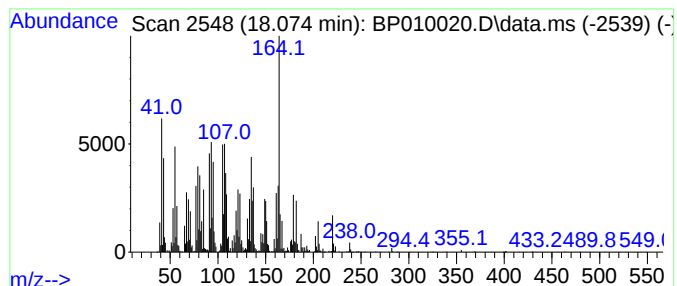
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (3S,3aS,6R,7R,9aS)-1,1,7-Tr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.074	4.28 ng/ul	396371	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(3S,3aS,6R,7R,9aS)-1,1,7-Trimeth...	238	C15H26O2	002649-64-1	99		
2	(3S,3aS,6R,7R,9aS)-1,1,7-Trimeth...	238	C15H26O2	002649-64-1	93		
3	2-Cyclopenten-1-one, 3-methyl-2-...	164	C11H16O	000488-10-8	38		
4	5-Pyrrolidin-2-ylidenemethyl-3,4-...	164	C9H12N2O	1010192-09-5	35		
5	Bicyclo[2.2.2]oct-2-ene, 1,2,3,6-...	164	C12H20	062376-14-1	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
Operator : CG/JU
Sample : N2373-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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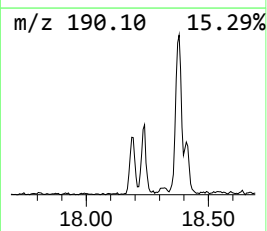
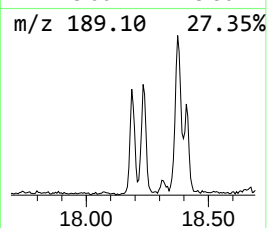
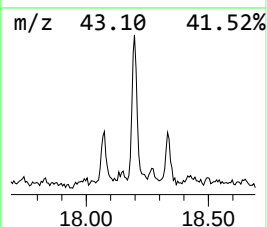
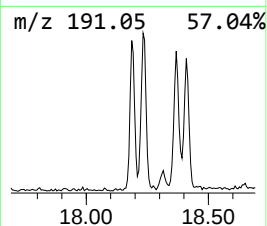
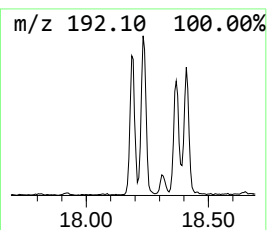
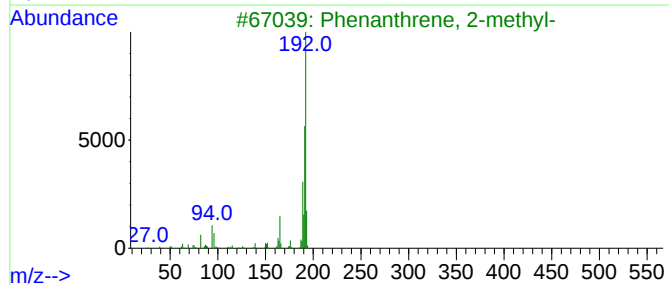
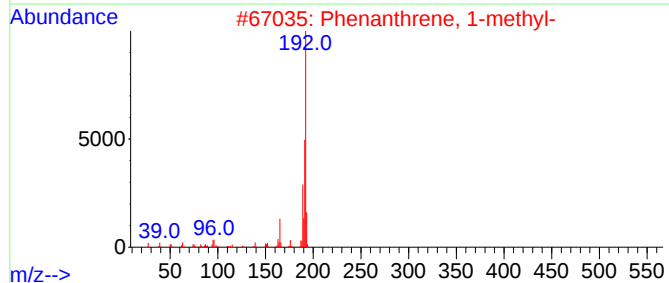
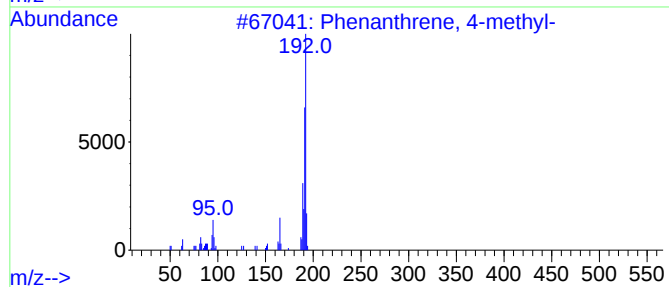
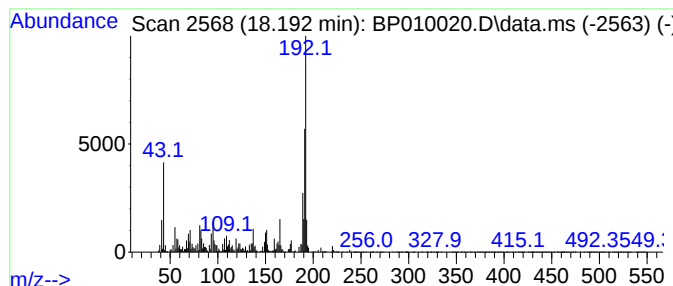
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	4.03 ng/ul	373581	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
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Misc :
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Instrument :
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ClientSampleId :
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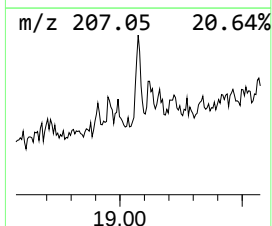
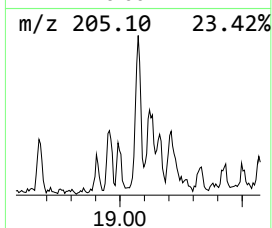
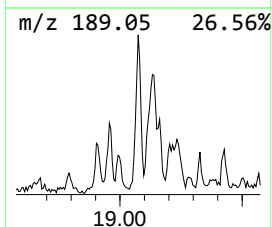
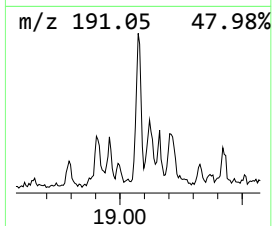
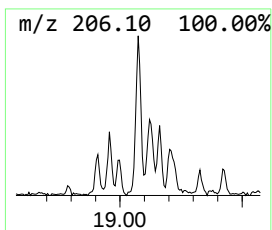
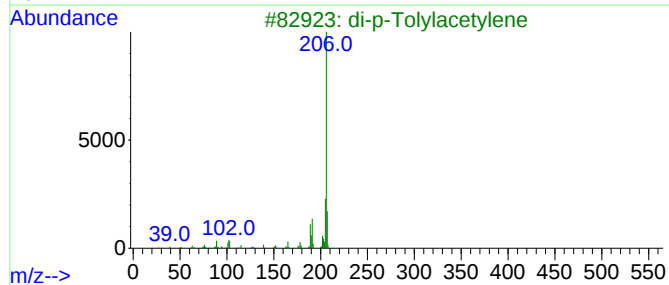
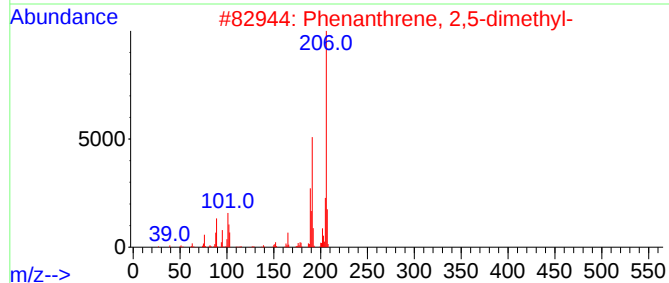
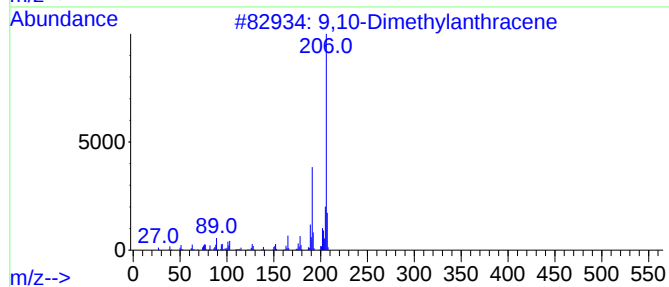
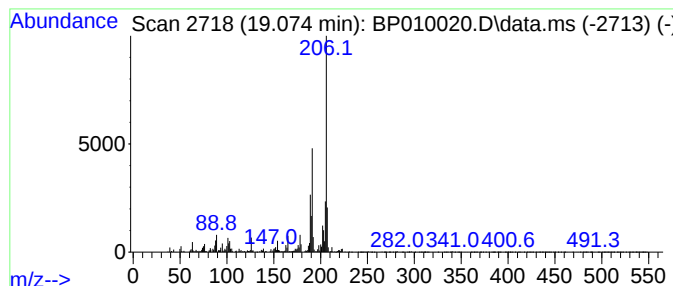
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9,10-Dimethylantracene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.074	2.32 ng/ul	214950	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
5			di-p-Tolylacetylene	206	C16H14	002789-88-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
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Sample : N2373-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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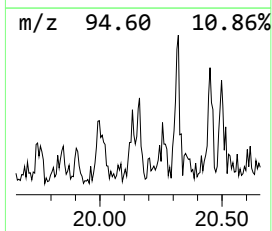
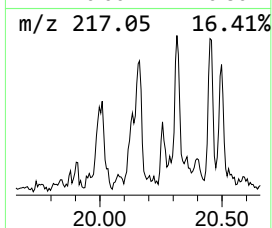
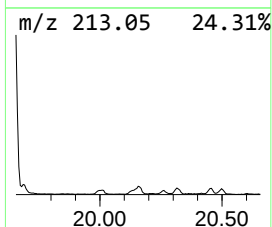
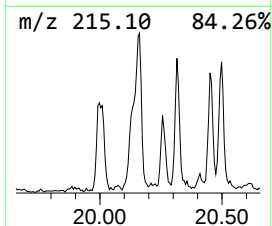
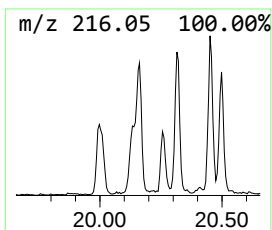
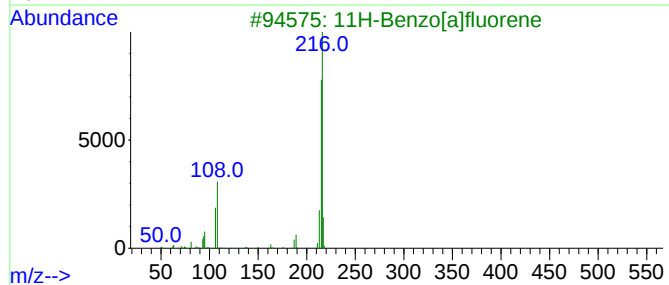
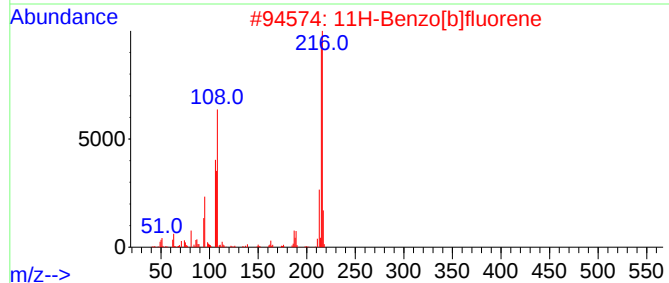
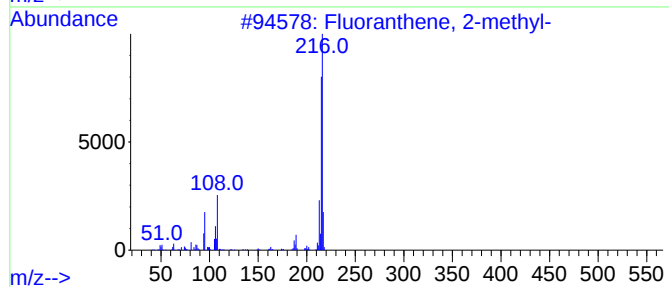
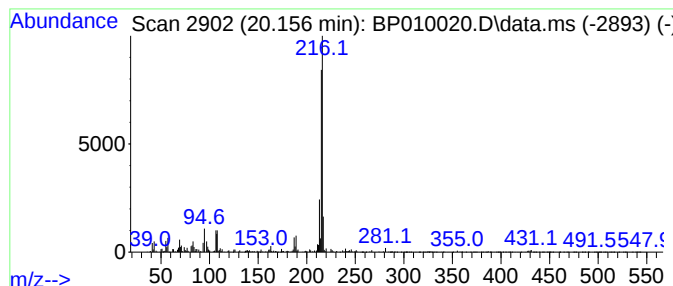
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Fluoranthene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.157	2.31 ng/ul	254863	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	86
5			Pyrene, 1-methyl-	216	C17H12	002381-21-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
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Misc :
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ClientSampleId :
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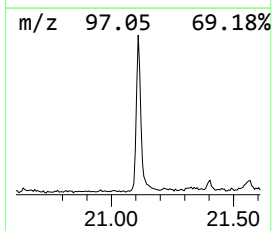
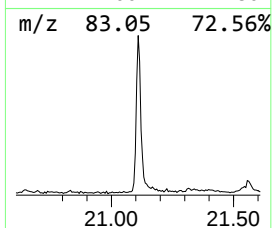
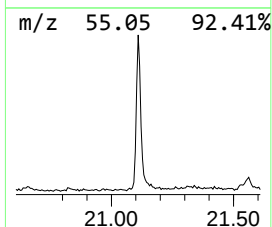
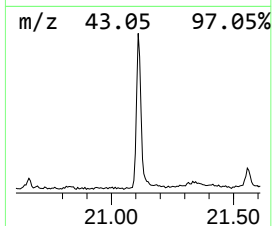
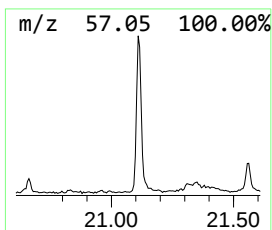
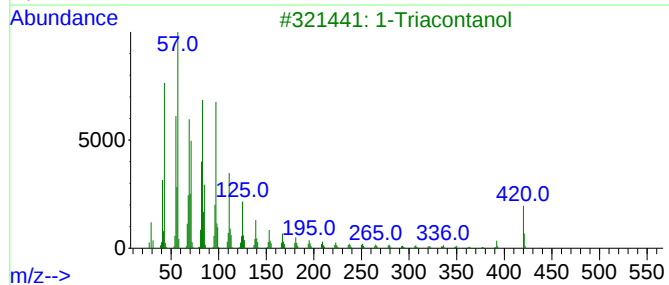
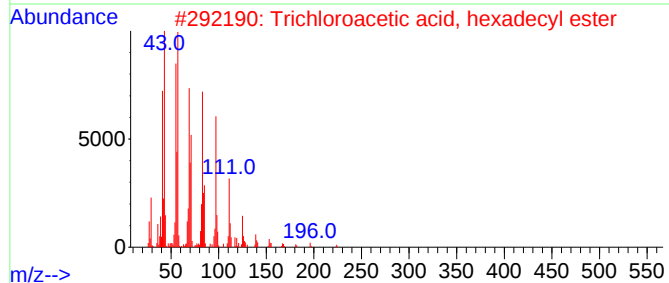
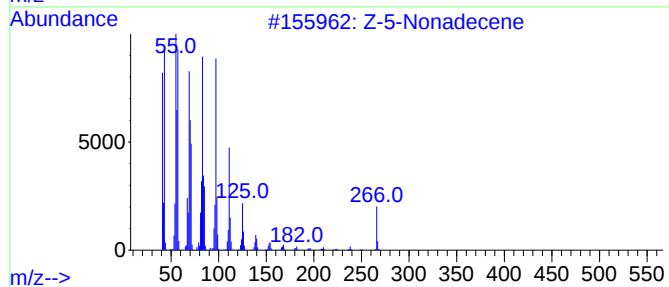
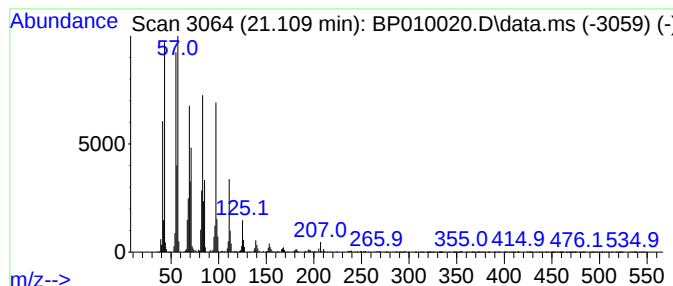
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.109	6.18 ng/ul	682856	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Z-5-Nonadecene	266	C19H38	1000131-11-8	95
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			1-Triacontanol	438	C30H62O	000593-50-0	94
4			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94
5			1-Docosene	308	C22H44	001599-67-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
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Misc :
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Instrument :
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ClientSampleId :
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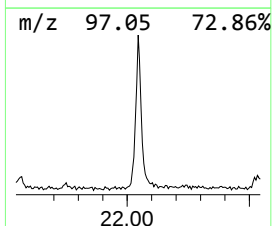
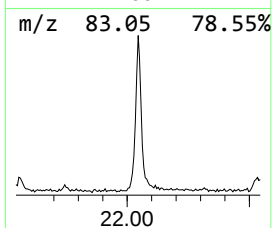
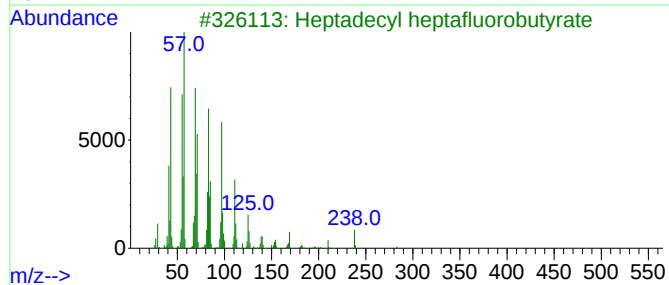
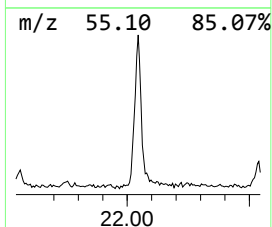
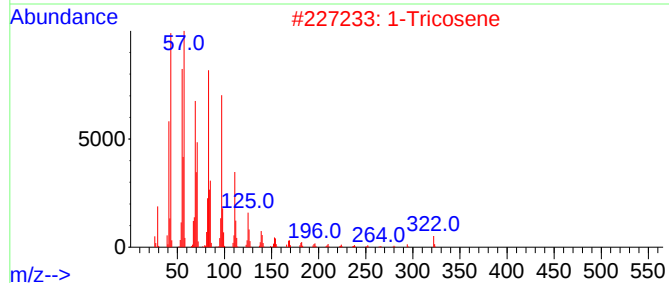
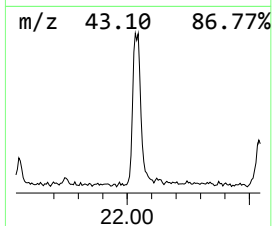
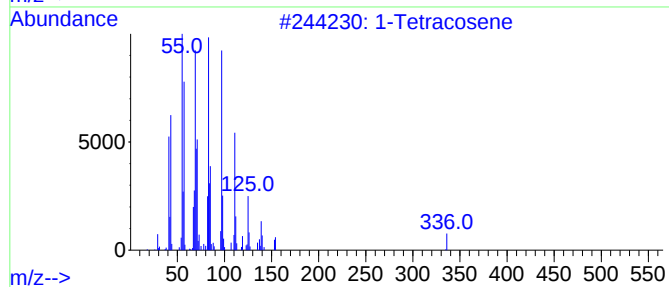
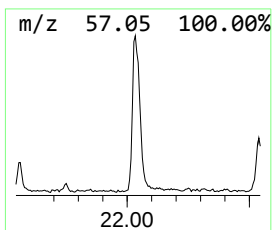
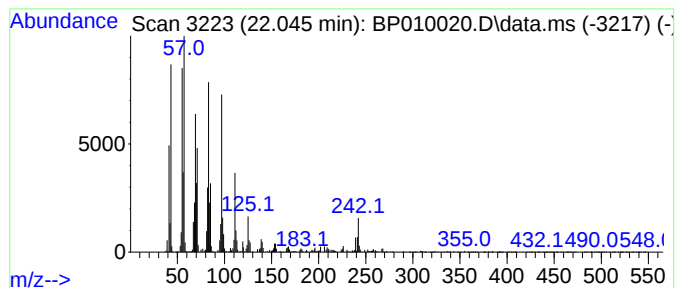
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	6.78 ng/ul	749443	Chrysene-d12	21.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	95
2		1-Tricosene	322	C23H46	018835-32-0	94
3		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94
5		Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
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ALS Vial : 15 Sample Multiplier: 1

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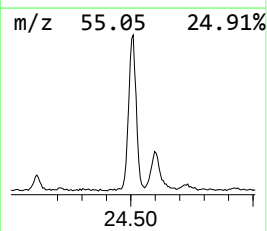
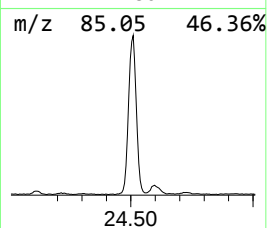
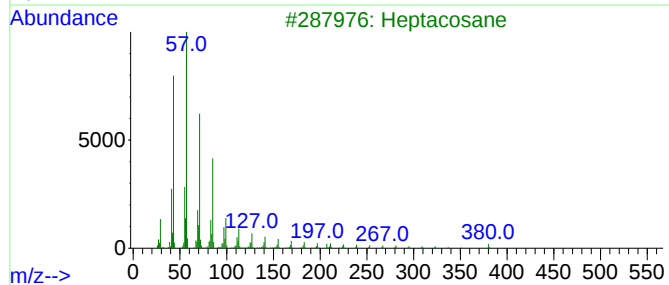
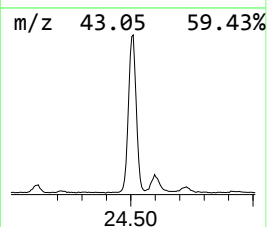
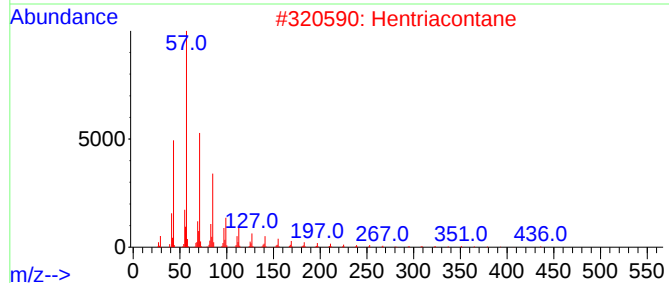
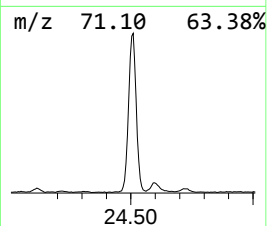
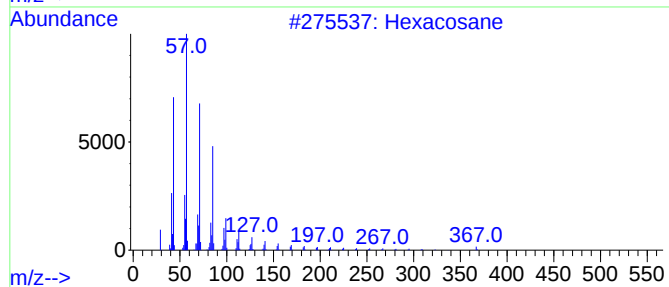
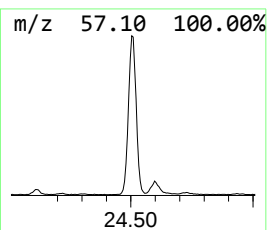
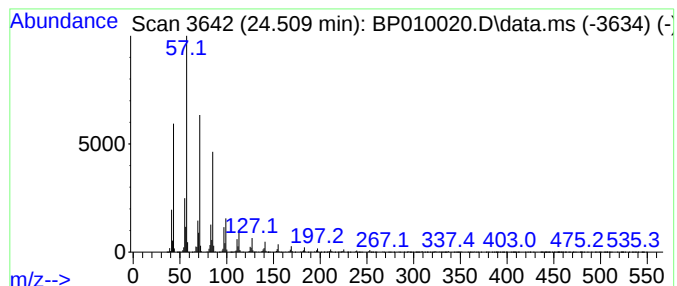
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.509	22.58 ng/ul	1964200	Perylene-d12	23.856

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	94
2		Hentriacontane	437	C31H64	000630-04-6	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Triacosane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
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Sample : N2373-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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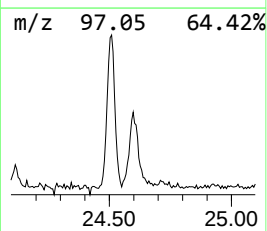
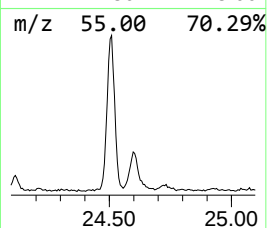
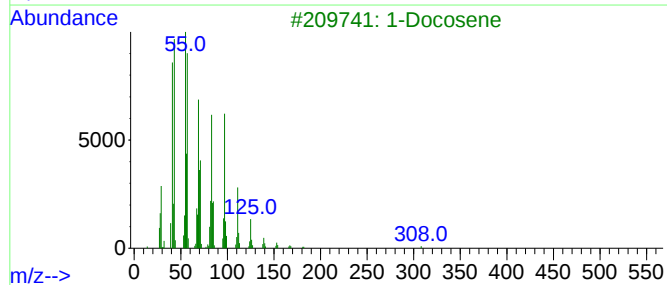
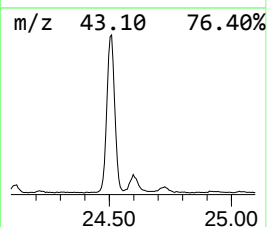
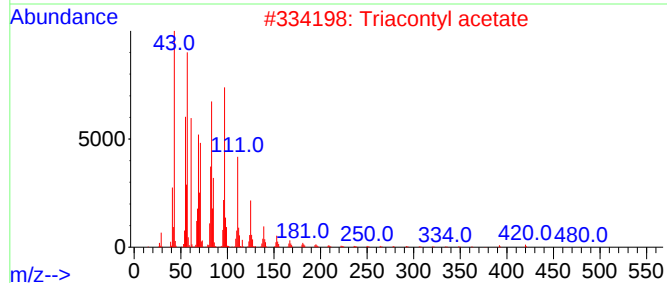
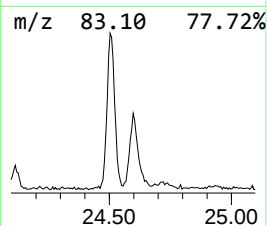
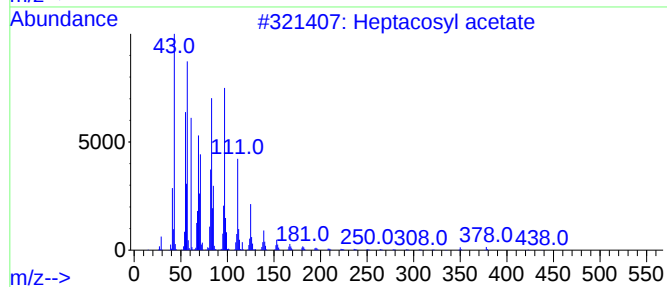
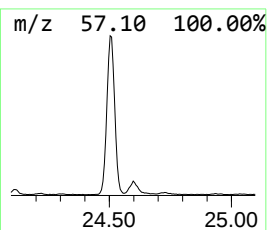
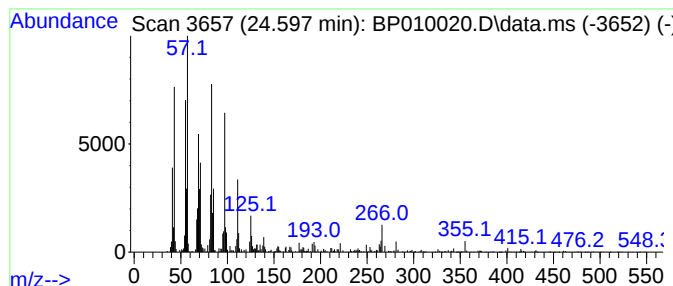
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Heptacosyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.597	3.38 ng/ul	294068	Perylene-d12	23.856

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosyl acetate	438	C29H58O2	1000351-78-2	93
2			Triacetyl acetate	480	C32H64O2	041755-58-2	92
3			1-Docosene	308	C22H44	001599-67-3	91
4			1-Heptacosanol	396	C27H56O	002004-39-9	91
5			Octacosanol	410	C28H58O	000557-61-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010020.D
Acq On : 21 Apr 2022 18:26
Operator : CG/JU
Sample : N2373-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFD2

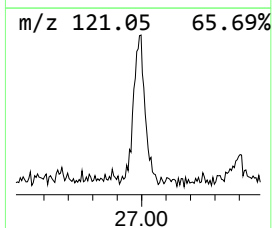
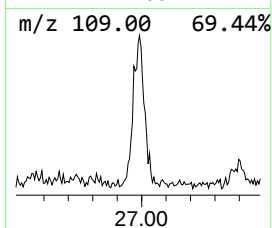
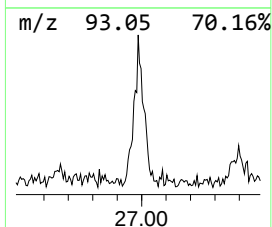
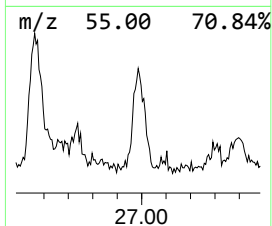
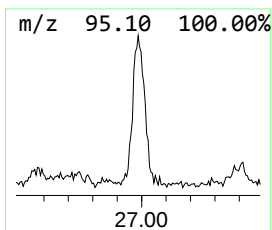
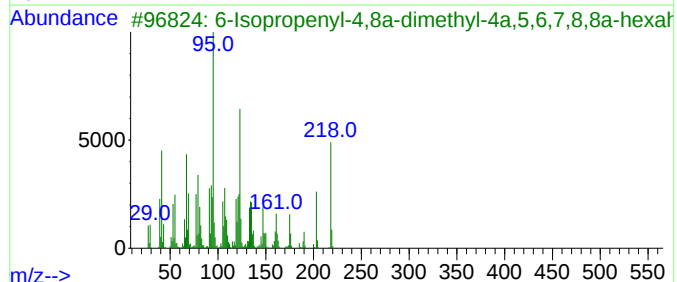
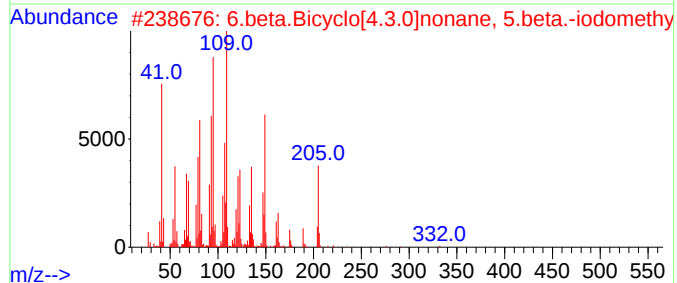
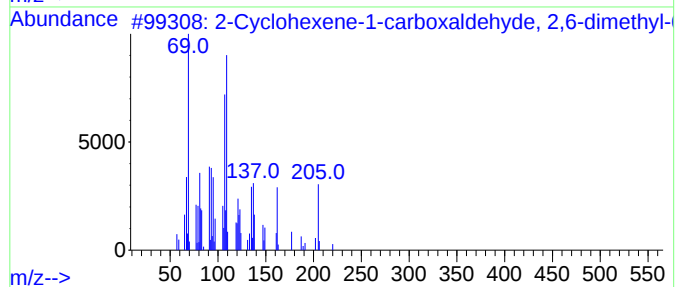
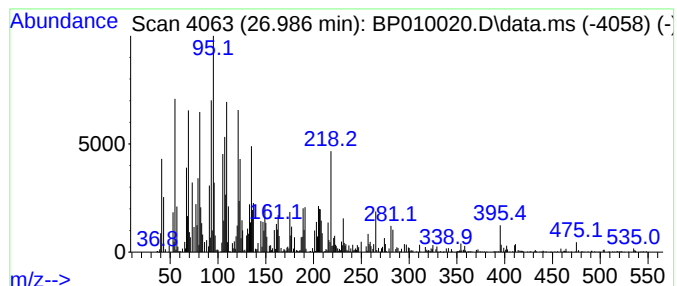
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.985	4.41 ng/ul	383554	Perylene-d12	23.856

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Cyclohexene-1-carboxaldehyde, ...	220	C15H24O	056772-07-7	42
2			6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	38
3			6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	38
4			(4aR,11aS)-4a-Methyl-3,4,4a,5,6...	218	C14H18O2	1000482-61-3	35
5			1,1-(Phenylthio)-2-isopropyl-cyc...	314	C19H22S2	122732-90-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010020.D
 Acq On : 21 Apr 2022 18:26
 Operator : CG/JU
 Sample : N2373-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFD2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.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
Eucalyptol	8.251	11.0	ng/ul	320238	1	7.939	584037	20.0
unknown-01	17.521	2.7	ng/ul	248503	4	17.321	1852200	20.0
(3S,3aS,6R,7R,9...)	18.074	4.3	ng/ul	396371	4	17.321	1852200	20.0
Phenanthrene, 4...	18.192	4.0	ng/ul	373581	4	17.321	1852200	20.0
9,10-Dimethylan...	19.074	2.3	ng/ul	214950	4	17.321	1852200	20.0
Fluoranthene, 2...	20.157	2.3	ng/ul	254863	5	21.404	2209520	20.0
(DEL) Alkane: S...	21.109	6.2	ng/ul	682856	5	21.404	2209520	20.0
(DEL) Alkane: S...	22.045	6.8	ng/ul	749443	5	21.404	2209520	20.0
(DEL) Alkane: S...	24.509	22.6	ng/ul	1964200	6	23.856	1739450	20.0
Heptacosyl acetate	24.597	3.4	ng/ul	294068	6	23.856	1739450	20.0
unknown-02	26.985	4.4	ng/ul	383554	6	23.856	1739450	20.0