

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
 Data File : BP022855.D
 Acq On : 05 Nov 2024 10:33
 Operator : RC/JU
 Sample : P4568-10
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4EH4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022855.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	29	32	41	rVB2	9919	13428	0.12%	0.036%
2	3.899	150	155	160	rBV2	7003	11395	0.10%	0.030%
3	4.446	243	248	253	rVB7	6944	11425	0.10%	0.031%
4	4.793	302	307	314	rBV2	7815	14441	0.13%	0.039%
5	5.099	354	359	366	rVB	22691	33326	0.30%	0.089%
6	5.652	449	453	464	rBV7	7109	18192	0.16%	0.049%
7	5.863	483	489	496	rVB2	23339	34702	0.31%	0.093%
8	5.999	506	512	518	rBV	20209	30457	0.28%	0.081%
9	6.499	593	597	601	rVB	12318	16961	0.15%	0.045%
10	6.828	641	653	670	rVV2	161338	373391	3.38%	0.999%
11	6.969	670	677	684	rVV	121084	190935	1.73%	0.511%
12	7.046	684	690	697	rVV	111409	179869	1.63%	0.481%
13	7.116	697	702	706	rVV	17516	26226	0.24%	0.070%
14	7.175	706	712	721	rVB	172371	272390	2.46%	0.729%
15	7.622	778	788	797	rBV	395761	616326	5.57%	1.649%
16	7.752	804	810	818	rVB	59830	103568	0.94%	0.277%
17	8.334	902	909	920	rBV	192273	348542	3.15%	0.932%
18	8.434	920	926	939	rVB	177278	319692	2.89%	0.855%
19	8.763	974	982	995	rBV	159878	285866	2.58%	0.765%
20	9.169	1045	1051	1058	rVB6	6957	13140	0.12%	0.035%
21	9.334	1073	1079	1086	rBV2	9721	20480	0.19%	0.055%
22	9.481	1095	1104	1118	rBV	143172	261340	2.36%	0.699%
23	9.693	1133	1140	1156	rBV2	40484	112236	1.01%	0.300%
24	9.840	1160	1165	1174	rVV2	20214	39437	0.36%	0.106%
25	10.022	1189	1196	1211	rBV	185269	411827	3.72%	1.102%
26	10.375	1248	1256	1263	rBV	542699	870036	7.87%	2.328%
27	10.540	1278	1284	1301	rBV3	37971	100288	0.91%	0.268%
28	10.781	1320	1325	1332	rBV3	5012	12707	0.11%	0.034%
29	10.940	1344	1352	1362	rBV3	10782	34717	0.31%	0.093%
30	11.204	1391	1397	1402	rBV4	6705	14378	0.13%	0.038%
31	11.775	1487	1494	1505	rBV4	6454	19553	0.18%	0.052%
32	12.051	1536	1541	1550	rVB4	6588	12402	0.11%	0.033%
33	12.269	1573	1578	1584	rVB4	6239	12391	0.11%	0.033%
34	13.557	1792	1797	1804	rBV10	6874	16555	0.15%	0.044%
35	13.657	1808	1814	1829	rVB	462907	661322	5.98%	1.769%
36	13.934	1852	1861	1879	rBV	538691	774883	7.01%	2.073%

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Integrator: RTE

Smoothing : OFF

Sampling : 1

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Stop Thrs : 0

Filtering: 5

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Max Peaks: 100

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
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37	14.245	1905	1914	1921	rBV	849314	1350729	12.21%	3.614%
38	14.310	1921	1925	1934	rVB	31307	59448	0.54%	0.159%
39	14.522	1953	1961	1980	rBV2	75082	236751	2.14%	0.633%
40	14.669	1980	1986	1995	rVV2	38255	85931	0.78%	0.230%

41	15.263	2080	2087	2093	rBV2	580940	1017319	9.20%	2.722%
42	15.322	2093	2097	2105	rVB	50737	74657	0.67%	0.200%
43	15.404	2106	2111	2120	rBV	89111	126940	1.15%	0.340%
44	15.516	2126	2130	2135	rVB8	8323	13893	0.13%	0.037%
45	15.645	2143	2152	2158	rBV	125632	200615	1.81%	0.537%

46	15.769	2169	2173	2180	rVB	17280	29592	0.27%	0.079%
47	15.986	2205	2210	2214	rBV5	12799	22619	0.20%	0.061%
48	16.175	2239	2242	2249	rVB9	5613	11560	0.10%	0.031%
49	16.351	2269	2272	2276	rVB4	11405	14033	0.13%	0.038%
50	16.445	2285	2288	2292	rBV6	8479	14180	0.13%	0.038%

51	16.710	2329	2333	2338	rBV3	14447	23866	0.22%	0.064%
52	16.857	2354	2358	2364	rVB	38048	55055	0.50%	0.147%
53	17.051	2384	2391	2395	rBV	1140251	1766456	15.97%	4.726%
54	17.098	2395	2399	2403	rVV	695543	924325	8.36%	2.473%
55	17.151	2403	2408	2412	rVV2	695912	1082591	9.79%	2.896%

56	17.186	2412	2414	2419	rVB	142660	171195	1.55%	0.458%
57	17.486	2459	2465	2468	rBV	48131	82267	0.74%	0.220%
58	17.751	2506	2510	2515	rBV8	12137	22945	0.21%	0.061%
59	17.851	2523	2527	2536	rBV5	21774	43561	0.39%	0.117%
60	17.951	2539	2544	2547	rVV	51330	78121	0.71%	0.209%

61	17.992	2547	2551	2560	rVB2	299242	445695	4.03%	1.192%
62	18.163	2575	2580	2584	rBV2	110041	182013	1.65%	0.487%
63	18.198	2584	2586	2595	rVB	39425	66287	0.60%	0.177%
64	18.486	2631	2635	2640	rVB	38877	62233	0.56%	0.166%
65	18.539	2640	2644	2647	rBV2	41159	68153	0.62%	0.182%

66	18.580	2648	2651	2654	rVV2	35754	42458	0.38%	0.114%
67	18.727	2669	2676	2682	rVB	7001858	11060440	100.00%	29.591%
68	18.816	2688	2691	2696	rVB	46481	67733	0.61%	0.181%
69	18.869	2697	2700	2703	rVB	37823	40289	0.36%	0.108%
70	18.922	2705	2709	2712	rBV2	32650	54670	0.49%	0.146%

71	19.186	2749	2754	2760	rVB	945519	1356138	12.26%	3.628%
72	19.245	2760	2764	2769	rVB	399612	493624	4.46%	1.321%
73	19.327	2774	2778	2784	rVB3	84946	143174	1.29%	0.383%
74	19.533	2807	2813	2816	rBV2	785409	1239500	11.21%	3.316%
75	19.569	2816	2819	2828	rVV	476038	701657	6.34%	1.877%

76	19.645	2829	2832	2838	rVB2	44094	70645	0.64%	0.189%
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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	19.751	2847	2850	2853	rBV	45731	56491	0.51%	0.151%
78	20.086	2903	2907	2911	rVB	98319	114844	1.04%	0.307%
79	20.186	2921	2924	2928	rVB2	89609	125272	1.13%	0.335%
80	20.969	3053	3057	3061	rBV5	105006	173780	1.57%	0.465%
81	21.069	3071	3074	3076	rBV	75643	99110	0.90%	0.265%
82	21.104	3077	3080	3085	rVB	315487	432671	3.91%	1.158%
83	21.486	3138	3145	3150	rBV2	1505061	2669563	24.14%	7.142%
84	21.527	3150	3152	3163	rVB	409411	557045	5.04%	1.490%
85	23.715	3517	3524	3532	rBV	233064	686611	6.21%	1.837%
86	24.515	3655	3660	3667	rVB2	217583	509599	4.61%	1.363%
87	24.739	3689	3698	3708	rBV	762888	2060965	18.63%	5.514%

Sum of corrected areas: 37378133

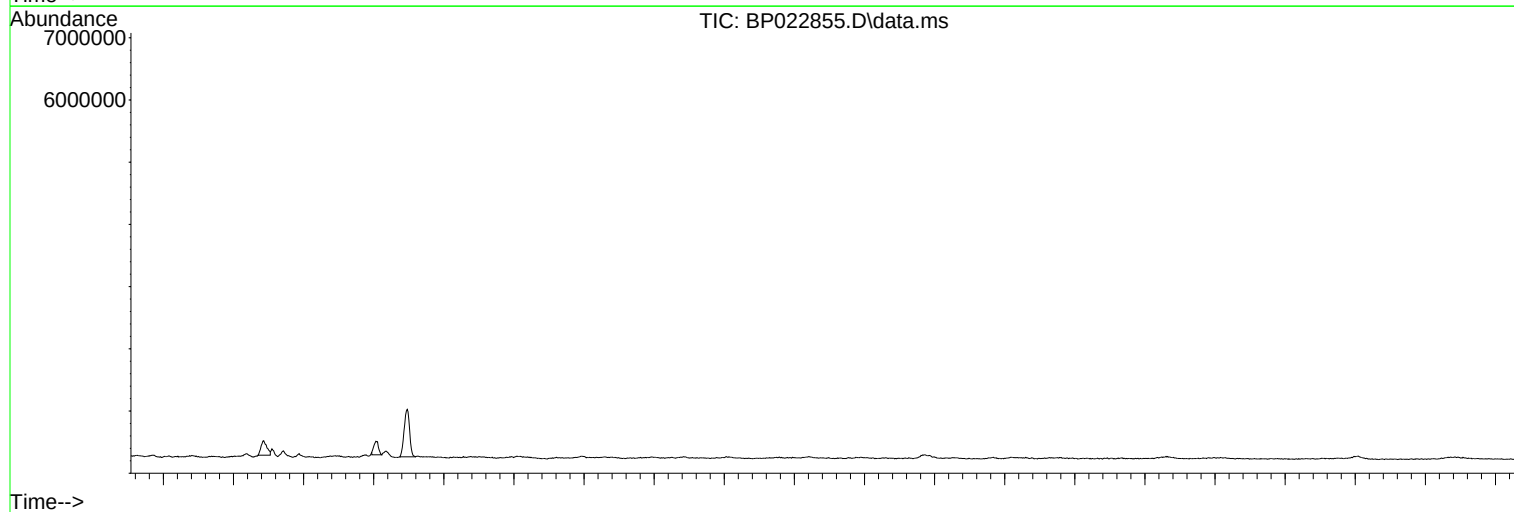
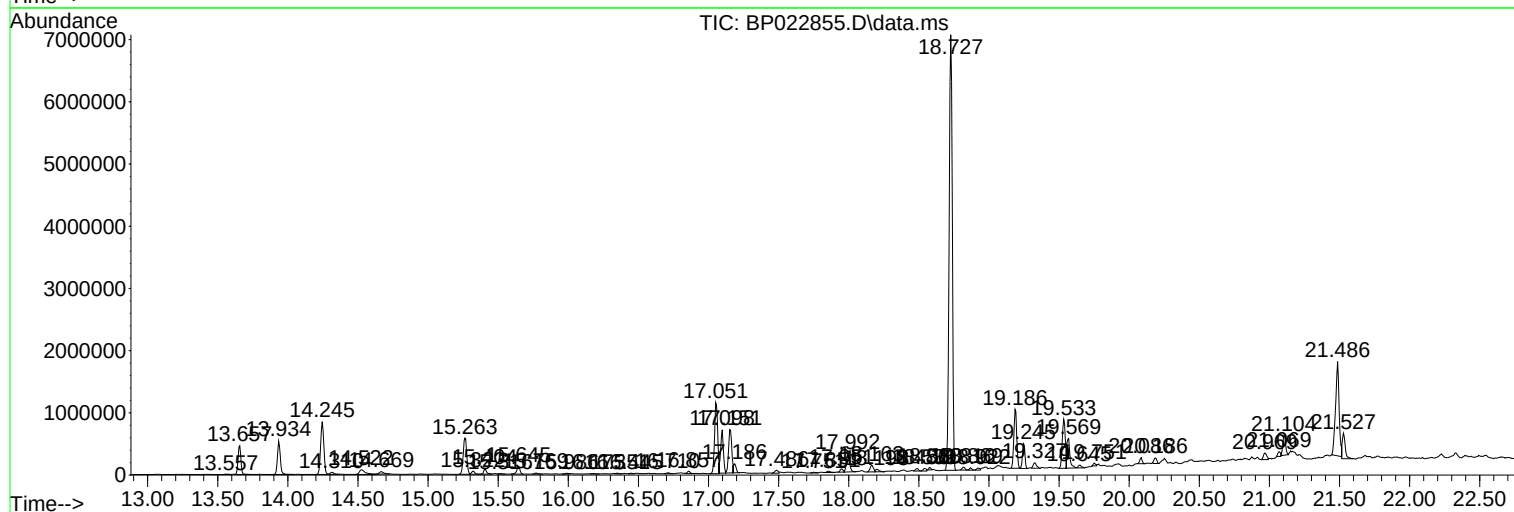
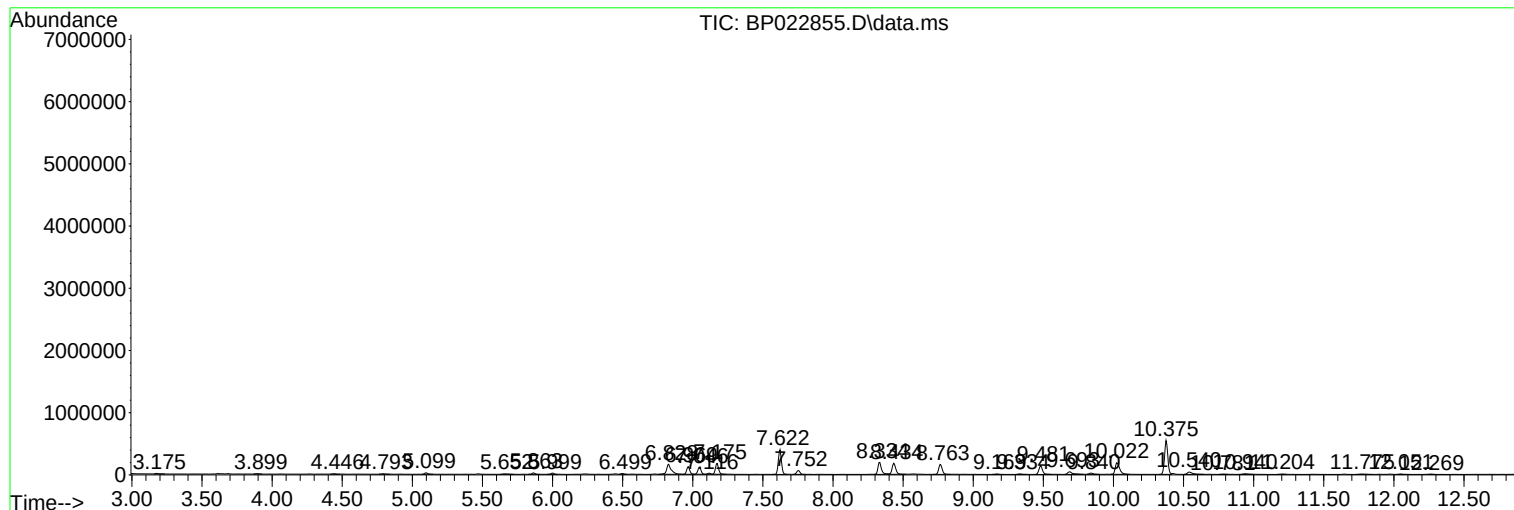
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P



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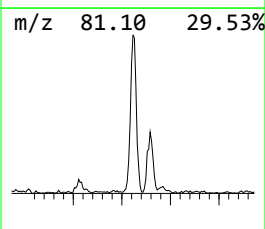
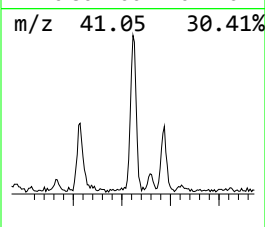
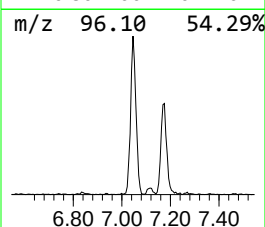
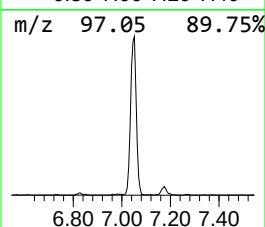
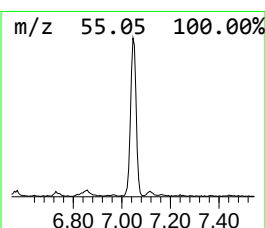
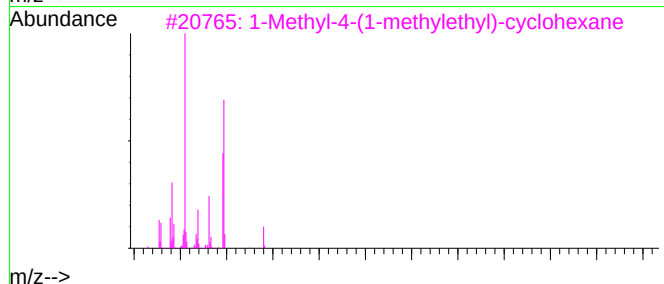
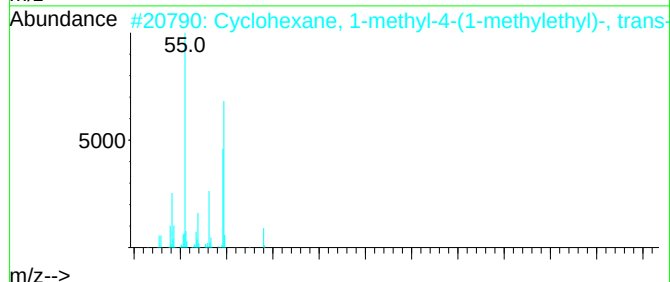
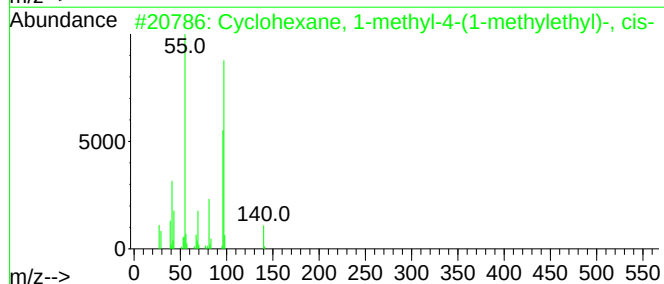
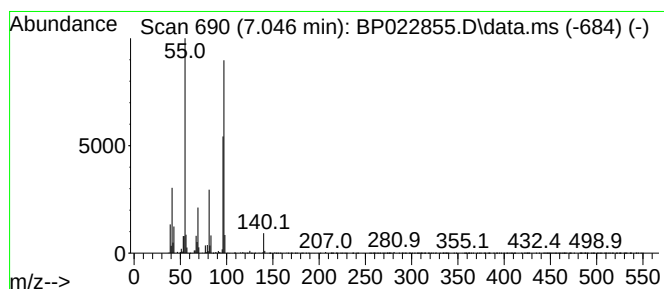
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Cyclic7.046 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.046	5.84 ng/ul	179869	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	97
2	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	97
3	1-Methyl-4-(1-methylethyl)-cyclo...	140	C10H20	000099-82-1	95
4	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	94
5	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	94



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

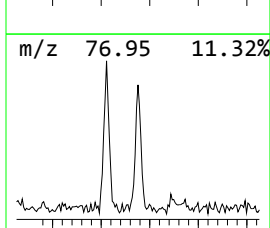
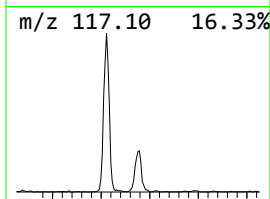
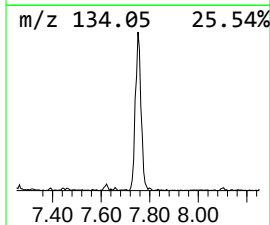
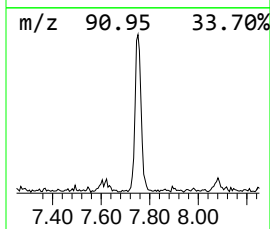
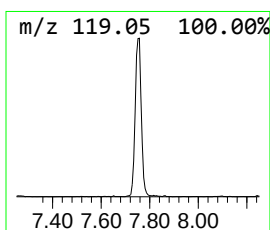
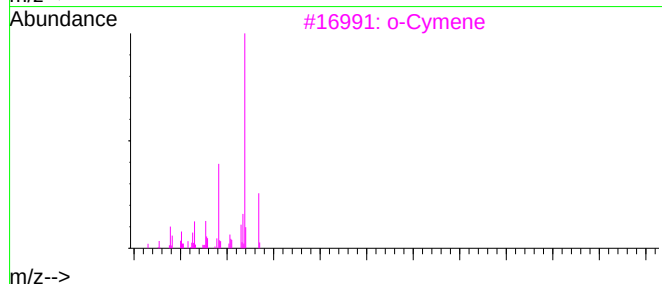
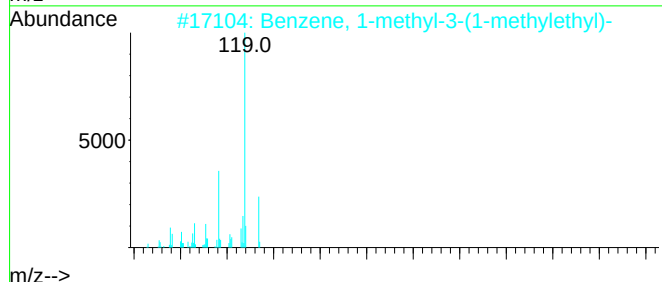
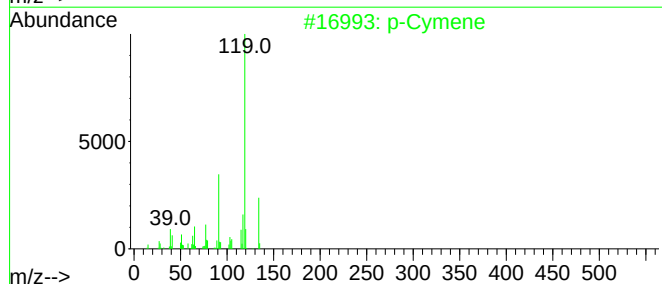
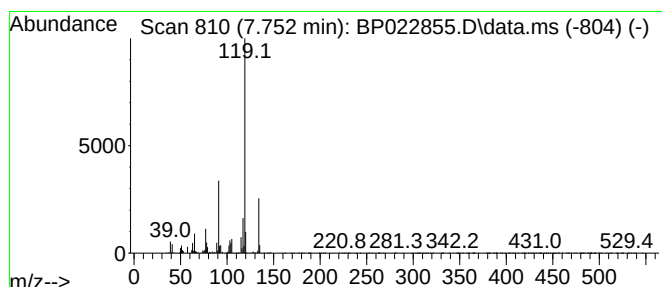
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 p-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.752	3.36 ng/ul	103568	1,4-Dichlorobenzene-d4	7.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	96
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	96
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5	p-Cymene	134	C10H14	000099-87-6	93



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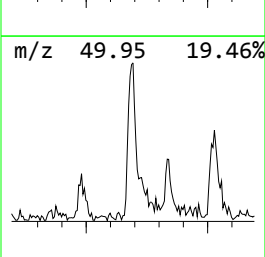
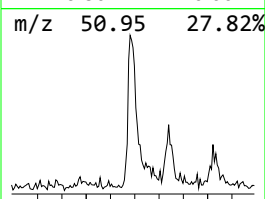
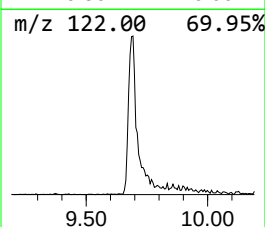
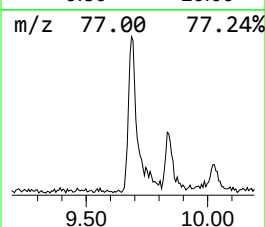
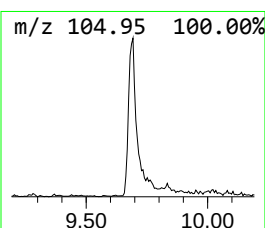
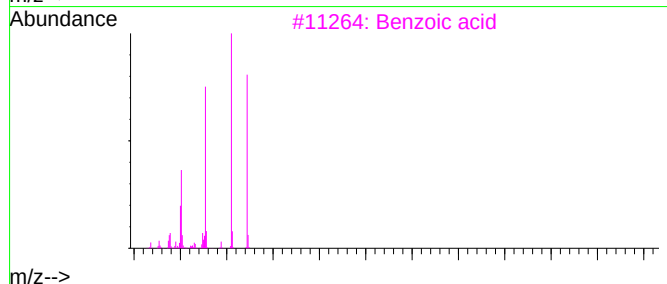
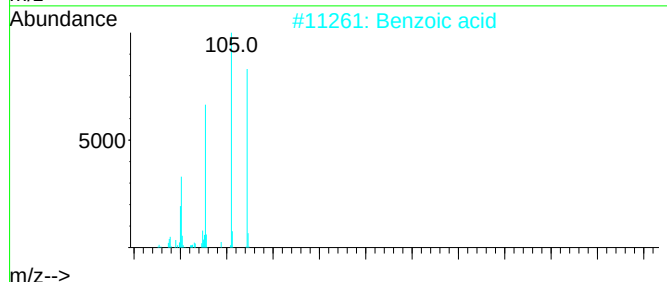
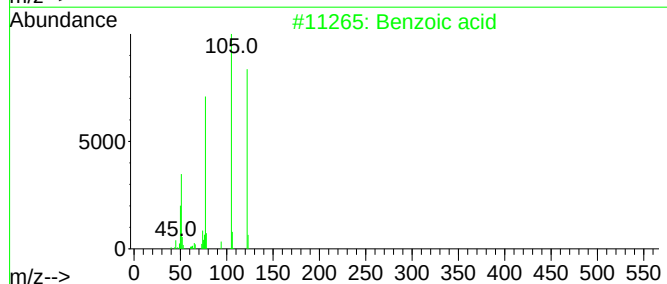
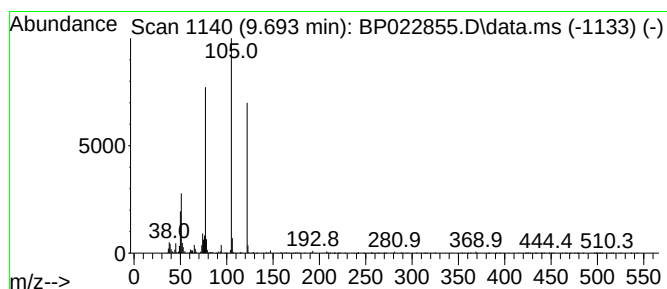
TIC Library : C:\DATABASE\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.693	2.58 ng/ul	112236	Naphthalene-d8	10.375

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid	122	C7H6O2	000065-85-0	94
2	Benzoic acid	122	C7H6O2	000065-85-0	94
3	Benzoic acid	122	C7H6O2	000065-85-0	91
4	.beta.-1,5-O-Dibenzoyl-ribofuranose	358	C19H18O7	1000129-71-8	89
5	Benzoic acid	122	C7H6O2	000065-85-0	86



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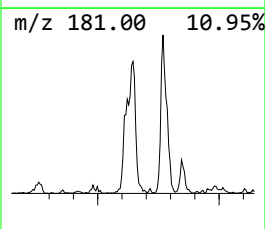
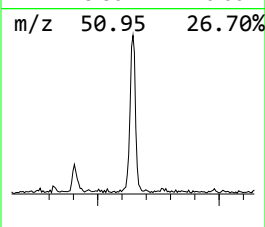
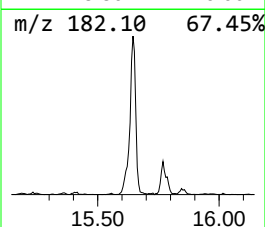
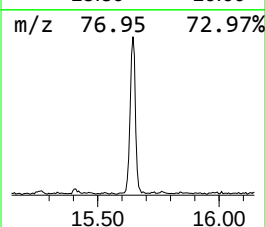
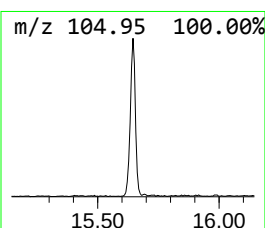
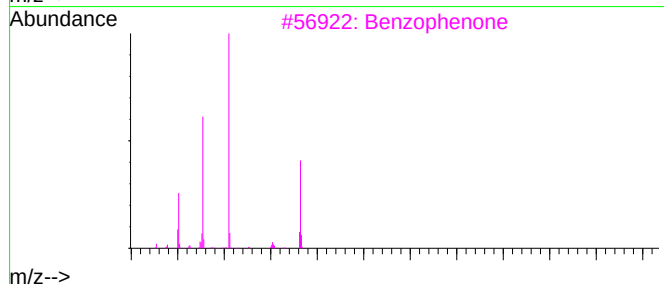
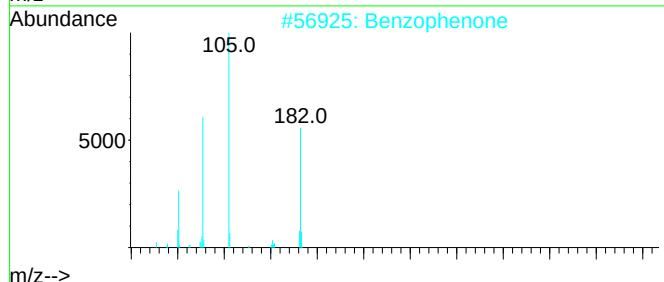
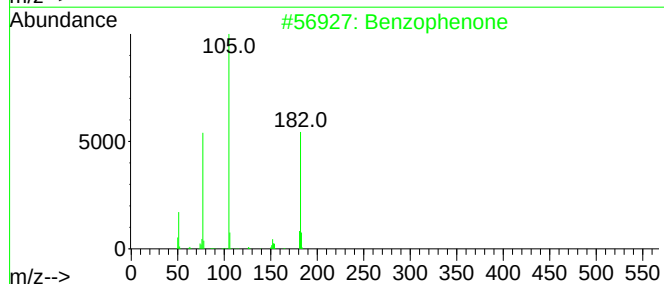
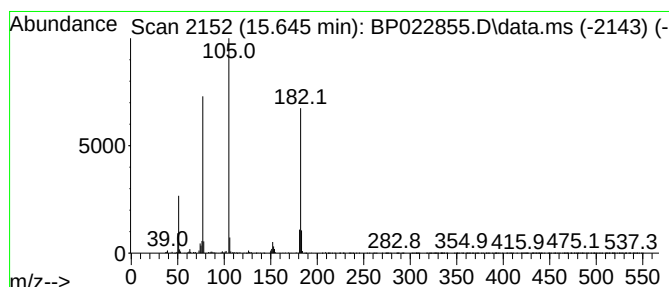
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.645	2.97 ng/ul	200615	Acenaphthene-d10	14.245	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	91
5	1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

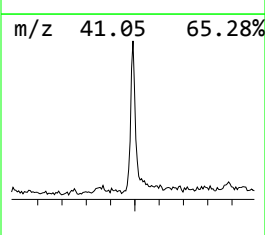
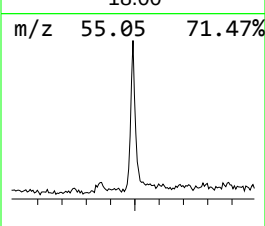
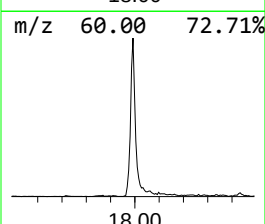
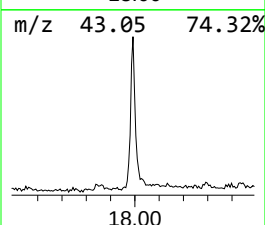
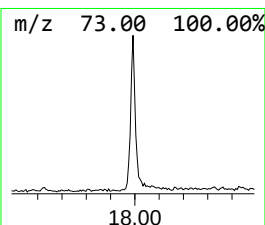
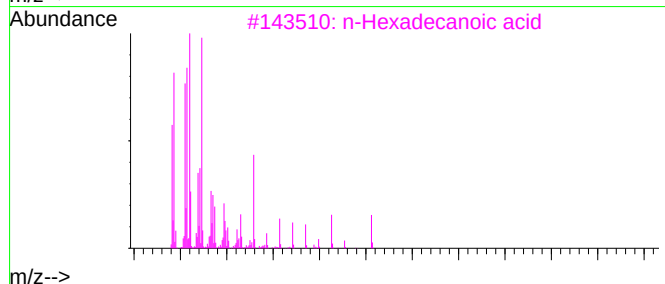
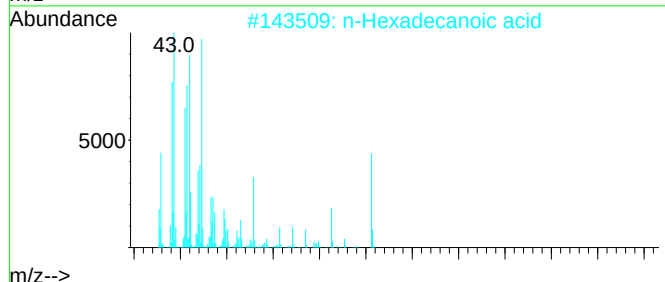
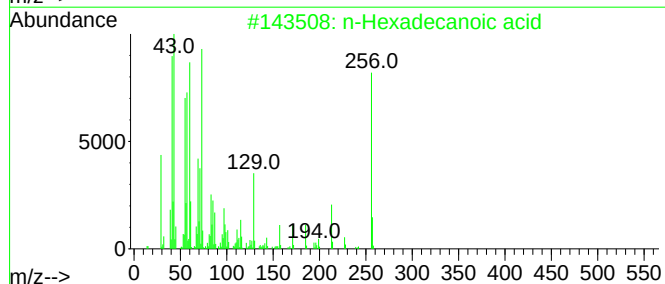
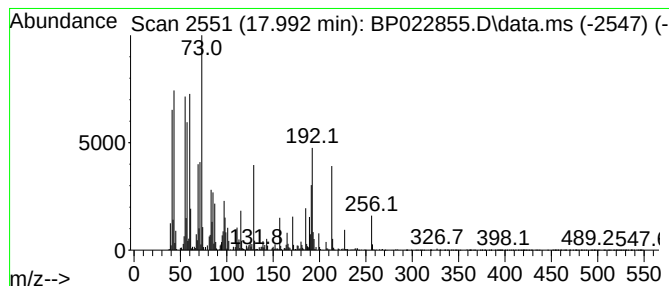
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.992	5.05 ng/ul	445695	Phenanthrene-d10	17.051	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
5	Tridecanoic acid	214	C13H26O2	000638-53-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

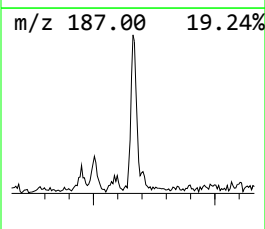
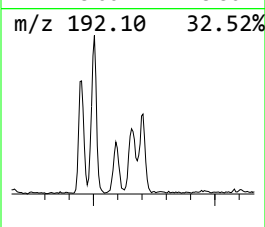
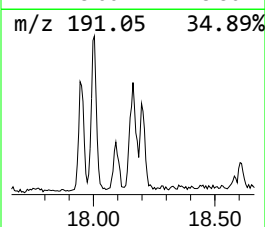
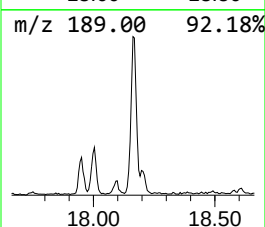
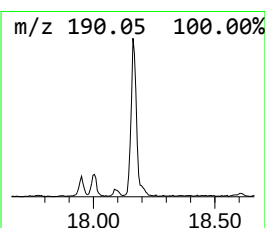
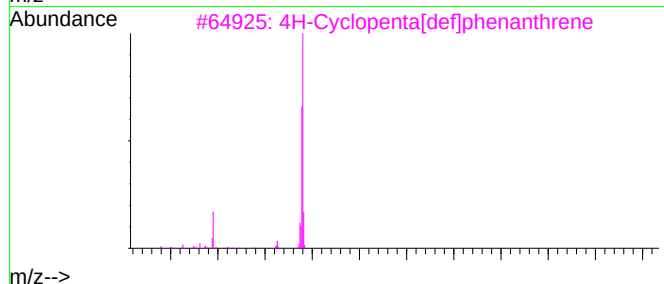
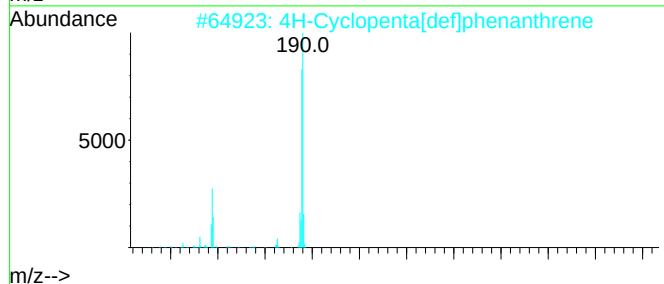
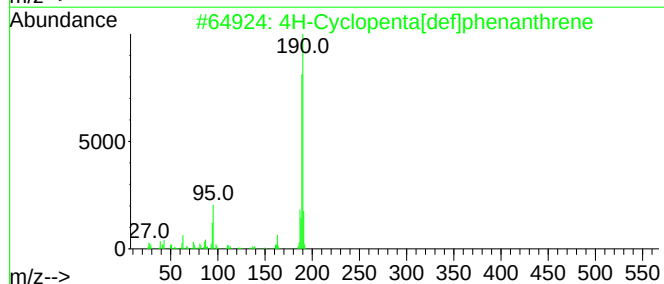
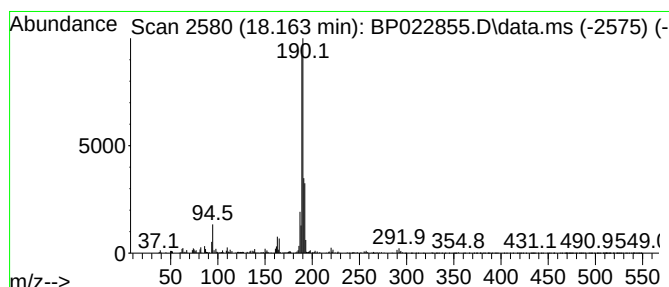
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.163	2.06 ng/ul	182013	Phenanthrene-d10	17.051	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	83
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	72
4	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BC103S	036155-87-0	59
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EH4

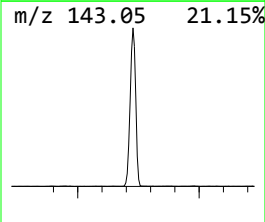
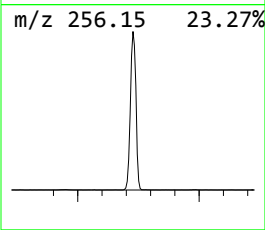
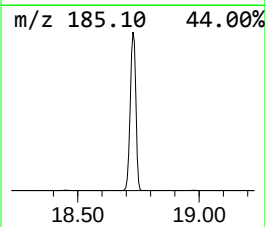
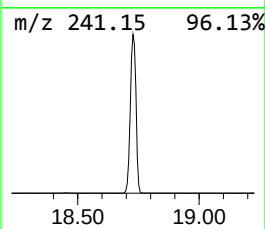
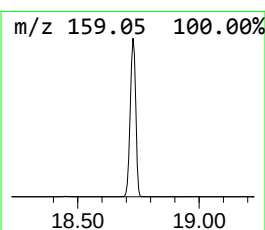
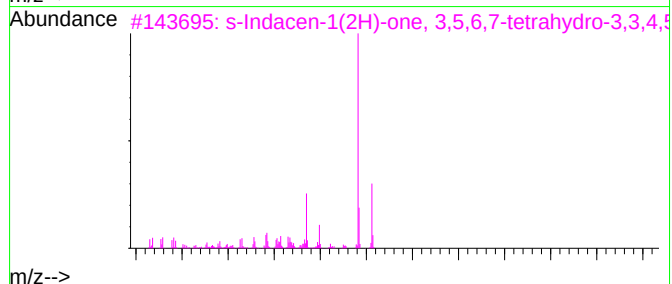
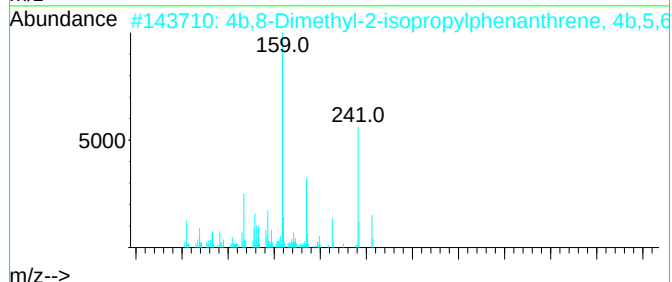
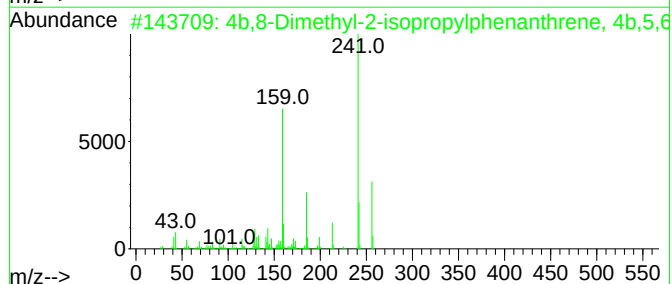
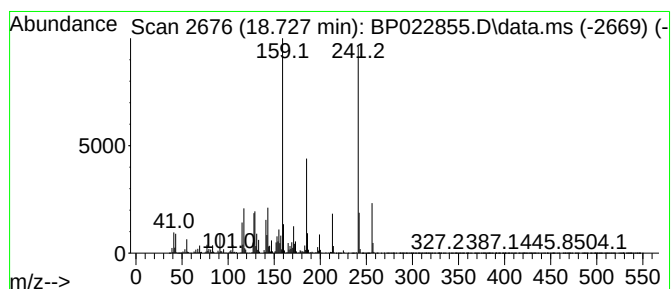
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4b,8-Dimethyl-2-isopropylph... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.727	125.23 ng/ul	11060400	Phenanthrene-d10	17.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	94
2			4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	83
3			s-Indacen-1(2H)-one, 3,5,6,7-tet...	256	C18H24O	038754-94-8	42
4			As-Indacene, 1,2,3,6,7,8-hexahyd...	256	C19H28	017465-47-3	30
5			Benzene, 1,1'-(1-methylethyliden...	256	C17H20O2	001568-83-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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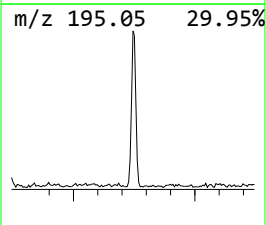
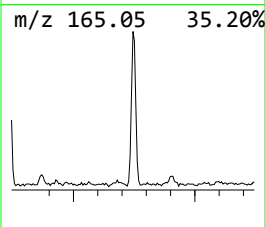
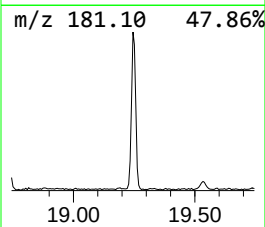
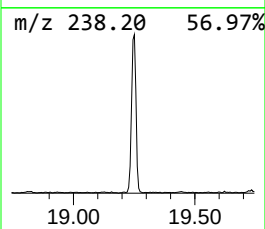
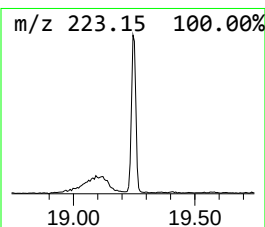
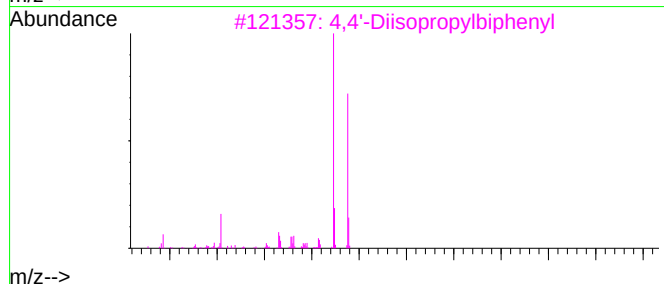
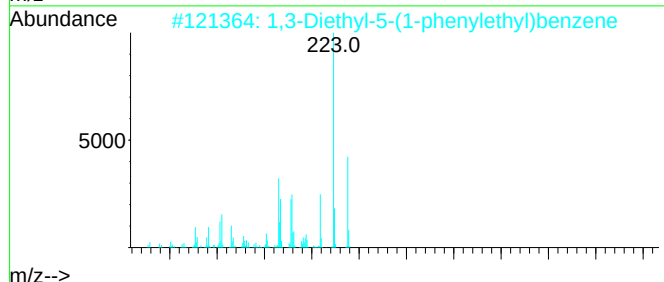
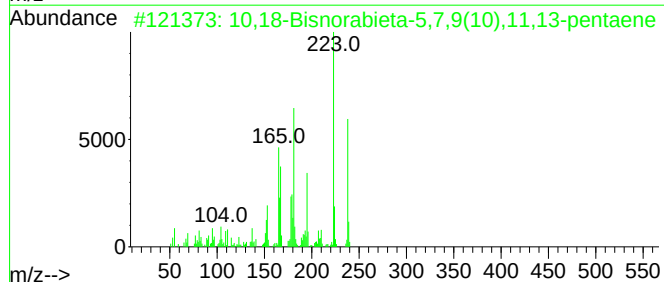
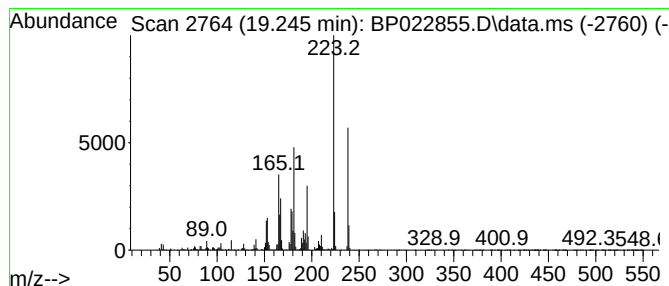
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.245	5.59 ng/ul	493624	Phenanthrene-d10	17.051

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2	1,3-Diethyl-5-(1-phenylethyl)ben...	238	C18H22	1000482-32-6	68
3	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
4	4-tert-Butylbenzenethiol, S-trim...	238	C13H22SSi	123045-43-2	49
5	Phenol, 2,6-bis(1,1-dimethylethy...	238	C14H22OS	000950-59-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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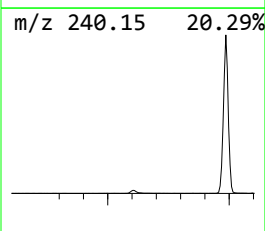
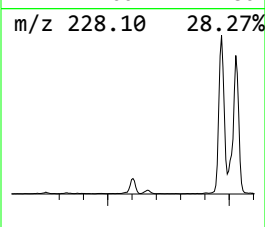
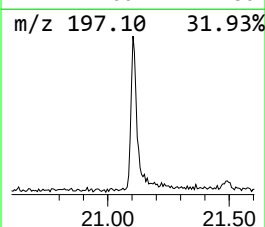
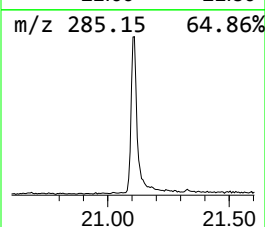
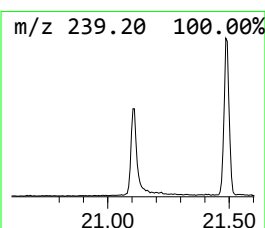
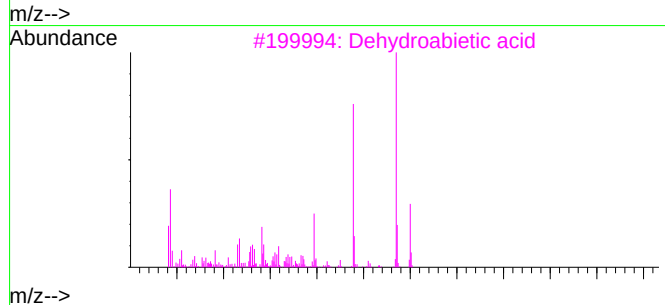
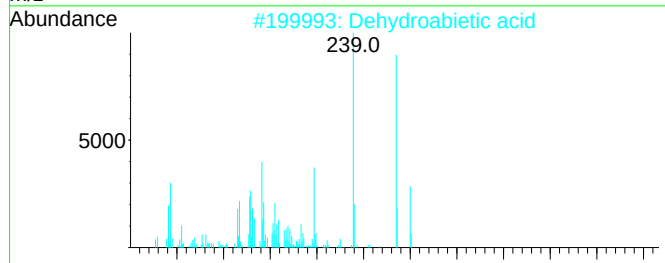
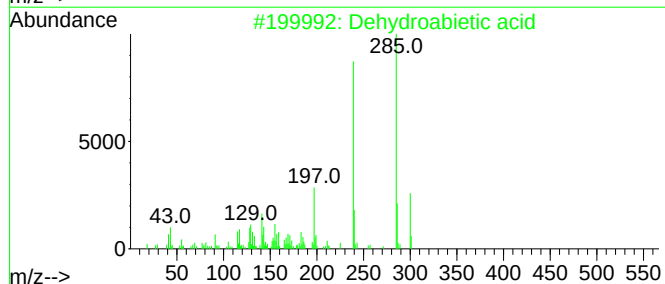
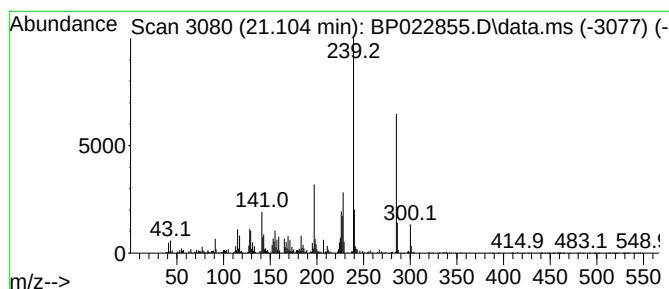
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Dehydroabietic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.104	3.24 ng/ul	432671	Chrysene-d12	21.486

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dehydroabietic acid	300	C20H28O2	001740-19-8	95
2	Dehydroabietic acid	300	C20H28O2	001740-19-8	93
3	Dehydroabietic acid	300	C20H28O2	001740-19-8	90
4	Dehydroabietic acid	300	C20H28O2	001740-19-8	90
5	1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022855.D
Acq On : 05 Nov 2024 10:33
Operator : RC/JU
Sample : P4568-10
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	7.046	5.8	ng/u1	179869	1	7.622	616326	20.0
p-Cymene	7.752	3.4	ng/u1	103568	1	7.622	616326	20.0
Benzoic acid	9.693	2.6	ng/u1	112236	2	10.375	870036	20.0
Benzophenone	15.645	3.0	ng/u1	200615	3	14.245	1350730	20.0
n-Hexadecanoic ...	17.992	5.0	ng/u1	445695	4	17.051	1766460	20.0
4H-Cyclopenta[d...	18.163	2.1	ng/u1	182013	4	17.051	1766460	20.0
4b,8-Dimethyl-2...	18.727	125.2	ng/u1	11060400	4	17.051	1766460	20.0
10,18-Bisnorabi...	19.245	5.6	ng/u1	493624	4	17.051	1766460	20.0
Dehydroabietic ...	21.104	3.2	ng/u1	432671	5	21.486	2669560	20.0