

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
 Data File : BP004659.D
 Acq On : 28 Jan 2021 15:30
 Operator : CG/JU
 Sample : M1203-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 X2Y84

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-BP012221.M
 Title : SVOA CALIBRATION

Signal : TIC: BP004659.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	10	13	16	rBV2	7248	9570	0.40%	0.022%
2	3.299	16	19	25	rVB	33021	42059	1.75%	0.097%
3	3.681	82	84	97	rBV	72234	112875	4.69%	0.260%
4	4.005	133	139	142	rBV4	5318	8016	0.33%	0.018%
5	4.322	189	193	200	rBV2	17632	28104	1.17%	0.065%
6	4.916	288	294	302	rVB	88844	122850	5.10%	0.283%
7	6.958	634	641	650	rVB	203170	323182	13.42%	0.743%
8	7.128	664	670	679	rBV	125539	194515	8.08%	0.447%
9	7.216	679	685	697	rVV	193429	287099	11.93%	0.660%
10	7.322	697	703	706	rVV	195495	334808	13.91%	0.770%
11	7.357	706	709	716	rVB	252070	360290	14.97%	0.829%
12	7.793	775	783	791	rBV	172741	267360	11.11%	0.615%
13	8.328	865	874	880	rBV2	127338	305050	12.67%	0.702%
14	8.375	880	882	889	rVB4	14194	19398	0.81%	0.045%
15	8.499	889	903	911	rBV	224583	363639	15.10%	0.836%
16	8.599	914	920	930	rBV	140175	212631	8.83%	0.489%
17	8.869	960	966	973	rBV	237822	381350	15.84%	0.877%
18	8.946	973	979	983	rVV	178620	307320	12.77%	0.707%
19	8.987	983	986	994	rVB	215939	345095	14.33%	0.794%
20	9.387	1044	1054	1059	rBV4	11087	18891	0.78%	0.043%
21	9.669	1092	1102	1109	rBV	168767	275064	11.43%	0.633%
22	10.004	1152	1159	1167	rBV	173031	265716	11.04%	0.611%
23	10.204	1185	1193	1195	rBV	285406	450860	18.73%	1.037%
24	10.228	1195	1197	1205	rVB	314537	471745	19.59%	1.085%
25	10.587	1245	1258	1262	rBV	220238	386714	16.06%	0.890%
26	10.640	1262	1267	1274	rVV	267127	443484	18.42%	1.020%
27	10.722	1274	1281	1293	rVB	223396	376174	15.63%	0.865%
28	10.928	1308	1316	1325	rVB	250225	407266	16.92%	0.937%
29	12.028	1499	1503	1510	rVB5	5466	7520	0.31%	0.017%
30	12.122	1512	1519	1525	rVV3	9660	15897	0.66%	0.037%
31	12.181	1525	1529	1536	rVB3	4406	6819	0.28%	0.016%
32	12.381	1557	1563	1568	rVV4	9014	15350	0.64%	0.035%
33	12.475	1568	1579	1589	rVB	392718	594068	24.68%	1.366%
34	12.610	1594	1602	1610	rBV3	339437	699348	29.05%	1.609%
35	12.863	1638	1645	1651	rBV	607164	875071	36.35%	2.013%
36	12.928	1651	1656	1663	rVV	358565	540001	22.43%	1.242%

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Integrator: RTE

Soothing : OFF

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Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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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-BP012221.M
 Title : SVOA CALIBRATION

37	13.275	1709	1715	1717	rVV	59806	85104	3.53%	0.196%
38	13.310	1717	1721	1735	rVB	421810	630402	26.19%	1.450%
39	13.557	1758	1763	1768	rBV8	4387	6780	0.28%	0.016%
40	13.851	1807	1813	1817	rVV	565609	793234	32.95%	1.825%
41	13.898	1817	1821	1831	rVV	628109	863377	35.86%	1.986%
42	14.016	1835	1841	1847	rVV	216719	304812	12.66%	0.701%
43	14.134	1850	1861	1863	rVV	528712	829169	34.44%	1.907%
44	14.157	1863	1865	1873	rVB	368490	494731	20.55%	1.138%
45	14.351	1893	1898	1903	rVB3	8656	12156	0.50%	0.028%
46	14.439	1906	1913	1919	rVV	379417	547321	22.73%	1.259%
47	14.504	1919	1924	1934	rVB	390566	562700	23.37%	1.294%
48	14.628	1937	1945	1953	rBV	293322	422729	17.56%	0.972%
49	14.716	1953	1960	1966	rVV	505564	703874	29.24%	1.619%
50	14.798	1969	1974	1977	rVV	196634	269605	11.20%	0.620%
51	14.839	1977	1981	1988	rVB	465067	639640	26.57%	1.471%
52	15.069	2017	2020	2024	rVB5	4168	6779	0.28%	0.016%
53	15.269	2046	2054	2058	rBV	550838	731528	30.39%	1.683%
54	15.304	2058	2060	2065	rVB	42434	50392	2.09%	0.116%
55	15.439	2076	2083	2087	rBV	753470	1079577	44.84%	2.483%
56	15.492	2087	2092	2097	rVV	871992	1183338	49.15%	2.722%
57	15.545	2097	2101	2111	rVB	272799	370206	15.38%	0.852%
58	15.675	2118	2123	2130	rBV2	9392	17689	0.73%	0.041%
59	16.092	2189	2194	2197	rBV7	4556	9153	0.38%	0.021%
60	16.128	2197	2200	2203	rVV4	6808	8280	0.34%	0.019%
61	16.169	2205	2207	2210	rVV4	7162	7885	0.33%	0.018%
62	16.222	2214	2216	2221	rVB6	5242	7807	0.32%	0.018%
63	16.339	2231	2236	2238	rBV5	4991	8988	0.37%	0.021%
64	16.386	2238	2244	2249	rVV	492140	631759	26.24%	1.453%
65	16.439	2249	2253	2256	rVV2	12611	17134	0.71%	0.039%
66	16.492	2257	2262	2270	rVV	356662	495519	20.58%	1.140%
67	16.581	2273	2277	2284	rBV7	12212	20507	0.85%	0.047%
68	16.839	2315	2321	2331	rBV	573607	802990	33.35%	1.847%
69	16.969	2339	2343	2349	rBV5	18355	26560	1.10%	0.061%
70	17.192	2374	2381	2387	rBV	498848	731147	30.37%	1.682%
71	17.298	2392	2399	2401	rBV	923737	1355651	56.31%	3.118%
72	17.328	2401	2404	2411	rVB	1146478	1513275	62.86%	3.481%
73	17.598	2443	2450	2460	rBV	936412	1297777	53.91%	2.985%
74	17.816	2483	2487	2491	rBV6	9427	12158	0.51%	0.028%
75	17.869	2492	2496	2501	rVB7	11098	17762	0.74%	0.041%
76	17.998	2513	2518	2520	rBV5	4642	7980	0.33%	0.018%

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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 >
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77	18.051	2525	2527	2530	rBV4	5748	7185	0.30%	0.017%
78	18.092	2530	2534	2539	rVV2	117928	151782	6.30%	0.349%
79	18.163	2541	2546	2551	rVB	857930	995657	41.36%	2.290%
80	18.386	2580	2584	2589	rBV7	7972	13476	0.56%	0.031%
81	18.580	2613	2617	2623	rBV2	33364	48615	2.02%	0.112%
82	18.745	2641	2645	2649	rVB3	17168	21918	0.91%	0.050%
83	18.833	2655	2660	2663	rBV	52431	64989	2.70%	0.149%
84	18.922	2671	2675	2681	rBV2	26140	38120	1.58%	0.088%
85	19.175	2715	2718	2722	rBV2	11193	14129	0.59%	0.032%
86	19.327	2740	2744	2752	rBV2	54552	78529	3.26%	0.181%
87	19.422	2757	2760	2762	rBV2	9883	12486	0.52%	0.029%
88	19.533	2773	2779	2785	rBV	1278550	1693386	70.34%	3.895%
89	20.439	2928	2933	2939	rBV	1098308	1173917	48.76%	2.700%
90	20.998	3024	3028	3034	rBV	260231	319577	13.27%	0.735%
91	21.198	3058	3062	3067	rBV	735272	753172	31.28%	1.732%
92	21.263	3068	3073	3081	rBV2	1410876	2407479	100.00%	5.538%
93	22.098	3210	3215	3221	rVB	1186559	1481358	61.53%	3.407%
94	22.892	3345	3350	3354	rBV	471895	873791	36.29%	2.010%
95	22.945	3354	3359	3365	rVB	1245740	2047312	85.04%	4.709%
96	23.451	3438	3445	3451	rBV	826436	1564312	64.98%	3.598%
97	23.510	3452	3455	3461	rVB2	95661	157382	6.54%	0.362%
98	23.598	3464	3470	3479	rVB	428820	793537	32.96%	1.825%
99	24.204	3570	3573	3580	rVB3	90181	154871	6.43%	0.356%
100	26.709	3991	3999	4011	rVB	369245	1154456	47.95%	2.656%

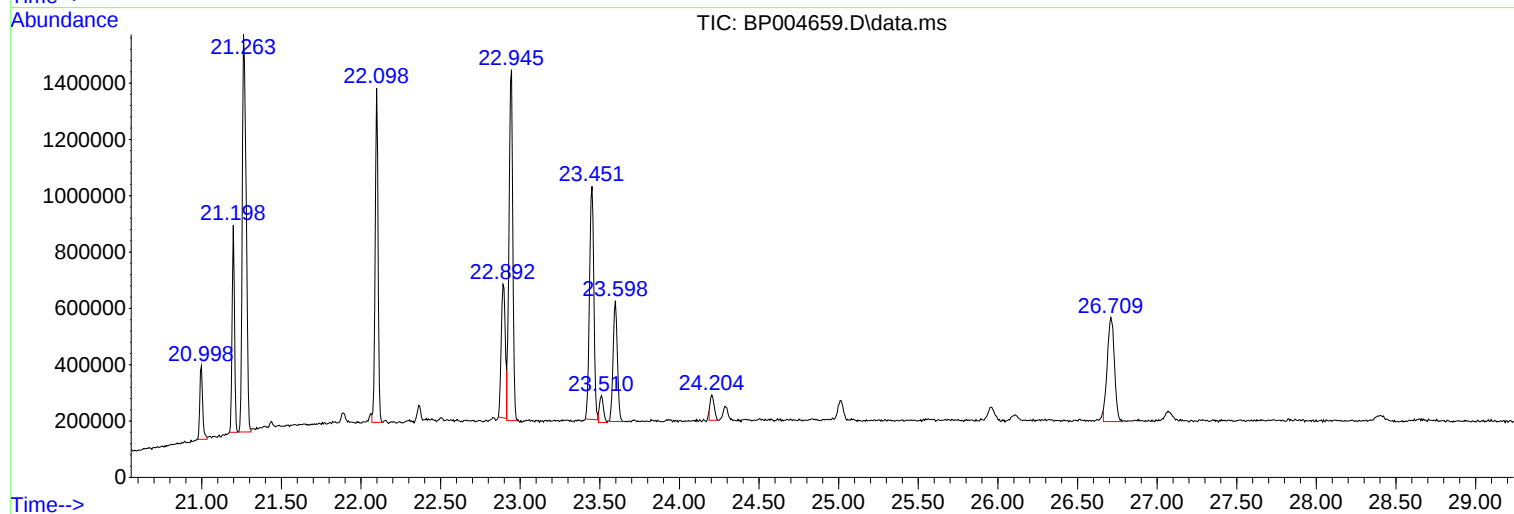
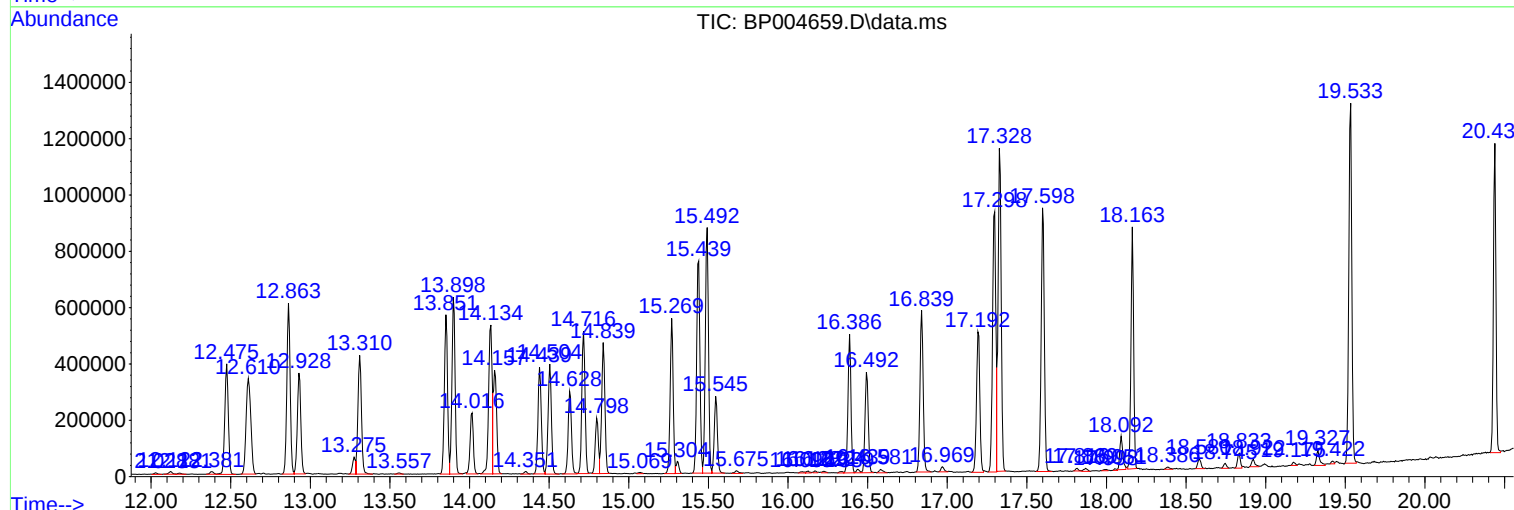
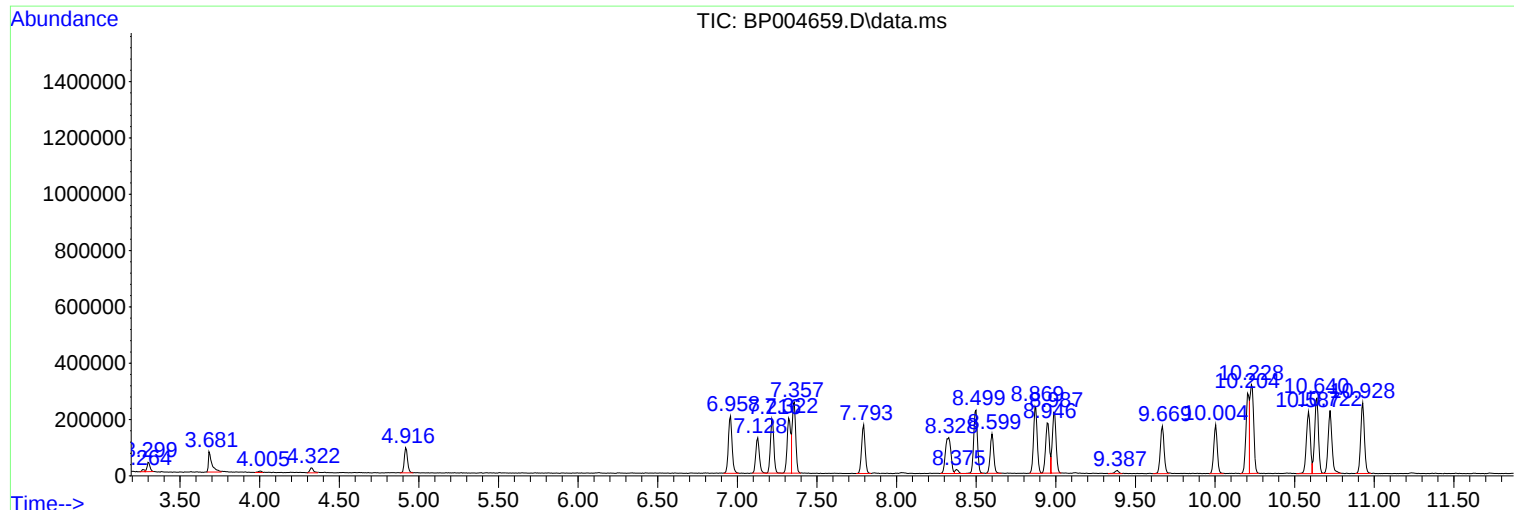
Sum of corrected areas: 43474110

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP012221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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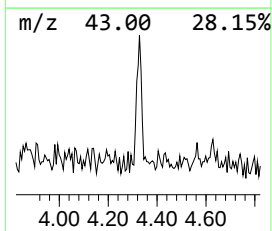
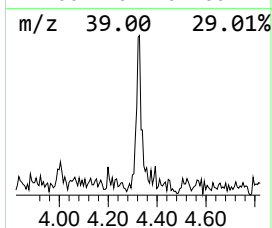
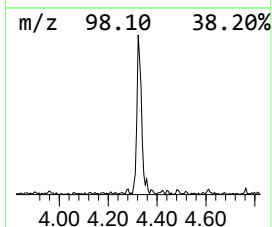
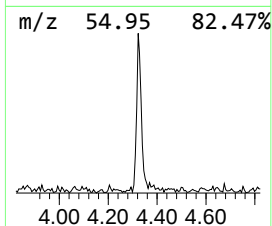
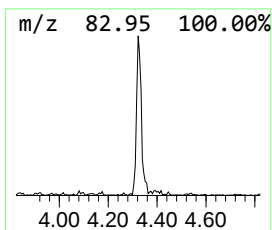
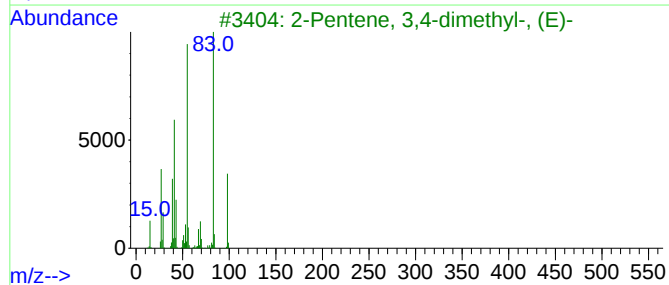
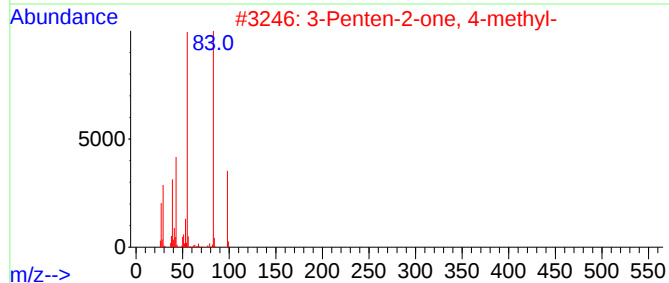
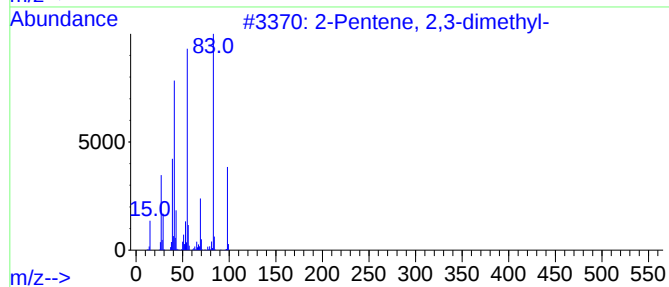
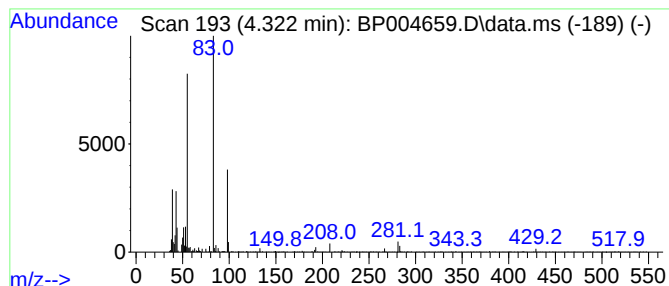
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP012221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.322	2.10 ng/ul	28104	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 2,3-dimethyl-			98	C7H14	010574-37-5	80
2	3-Penten-2-one, 4-methyl-			98	C6H10O	000141-79-7	80
3	2-Pentene, 3,4-dimethyl-, (E)-			98	C7H14	004914-92-5	80
4	2-Pentene, 3,4-dimethyl-, (Z)-			98	C7H14	004914-91-4	80
5	2-Pentene, 2,4-dimethyl-			98	C7H14	000625-65-0	72



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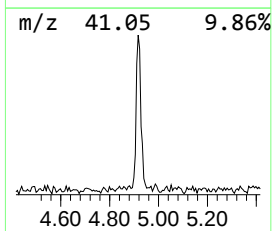
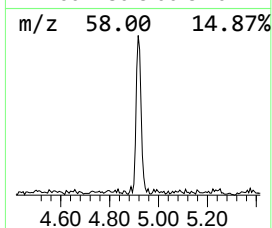
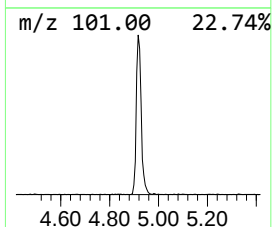
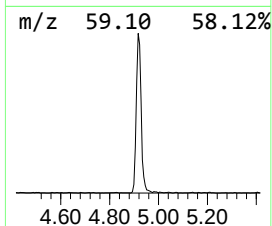
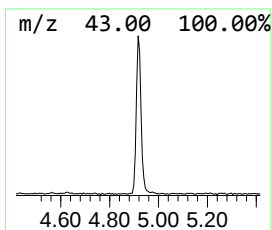
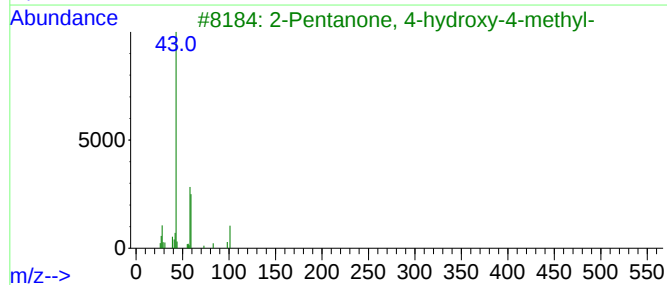
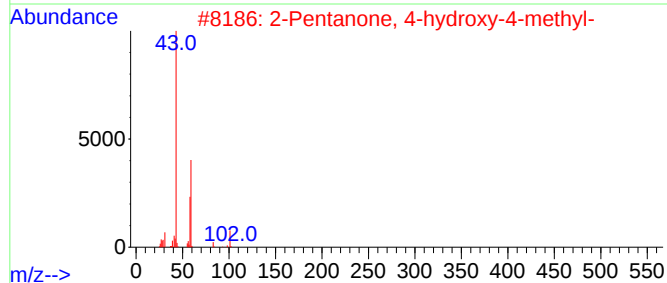
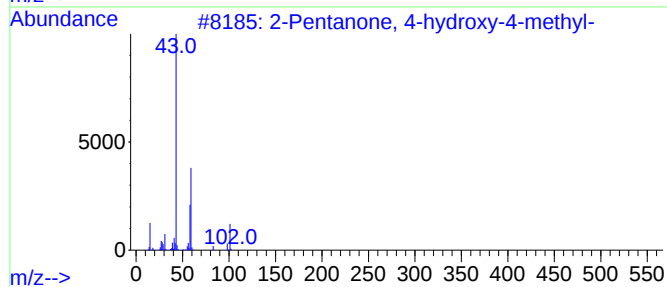
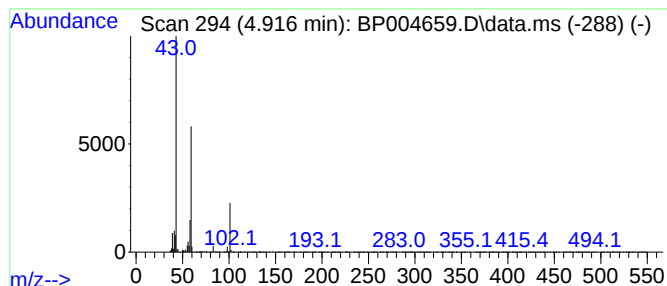
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.916	9.19 ng/ul	122850	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
5	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17		



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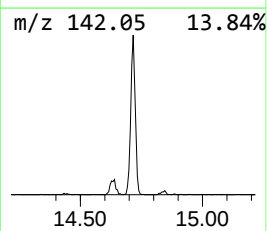
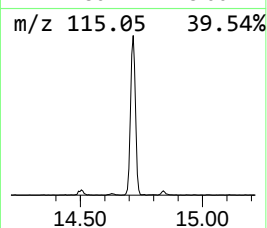
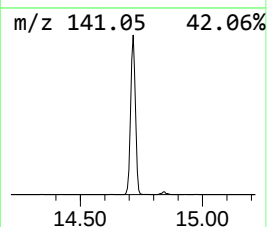
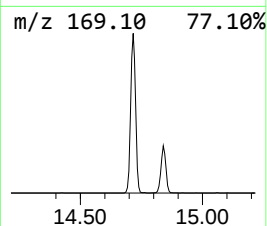
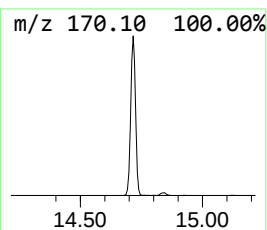
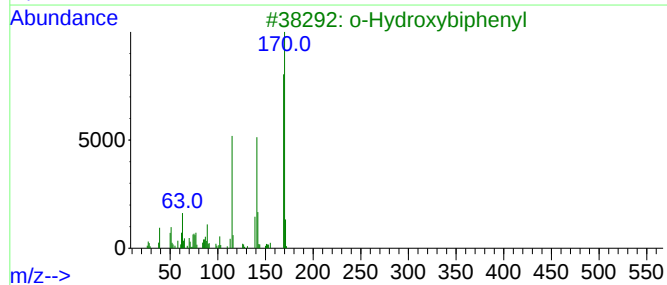
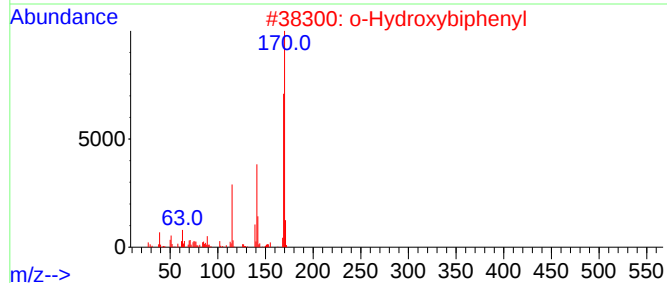
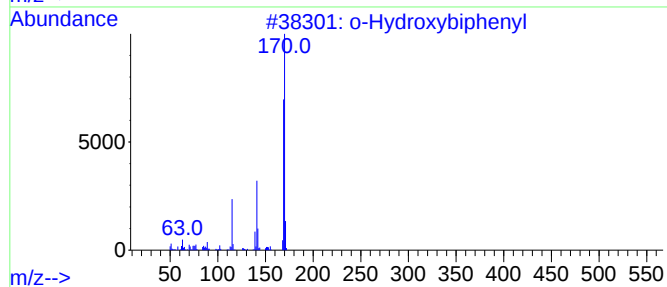
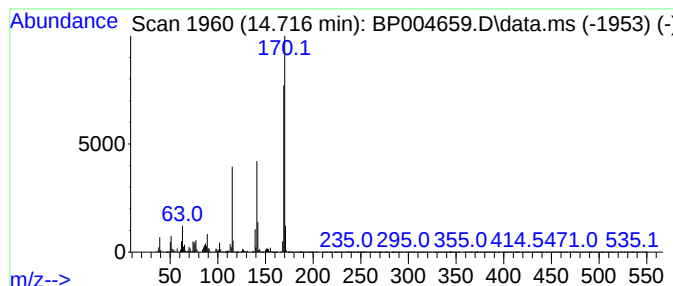
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 o-Hydroxybiphenyl Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.716	25.72 ng/ul	703874	Acenaphthene-d10	14.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	97
2			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	96
3			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	96
4			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	90
5			o-Hydroxybiphenyl	170	C12H10O	000090-43-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
Data File : BP004659.D
Acq On : 28 Jan 2021 15:30
Operator : CG/JU
Sample : M1203-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
X2Y84

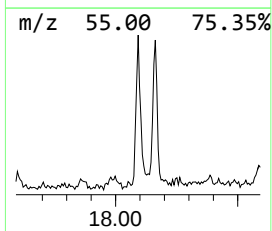
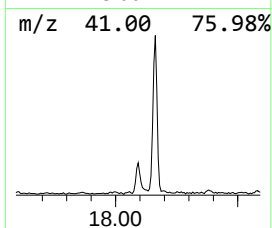
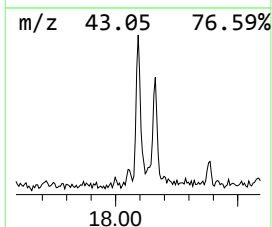
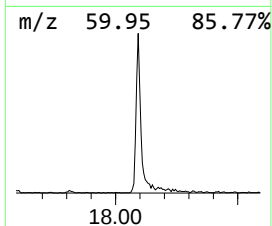
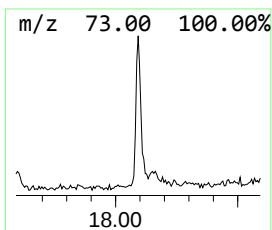
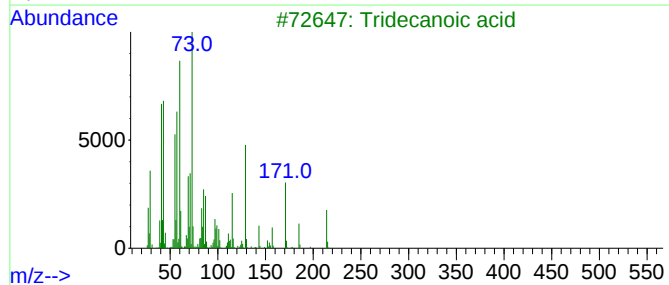
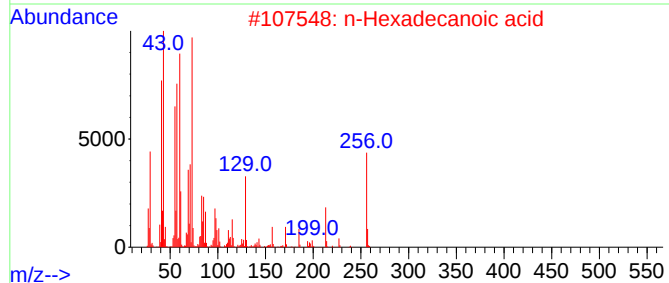
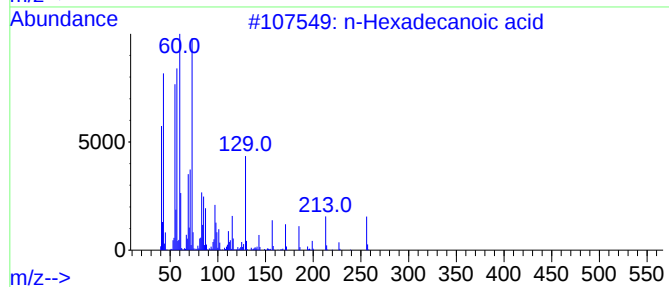
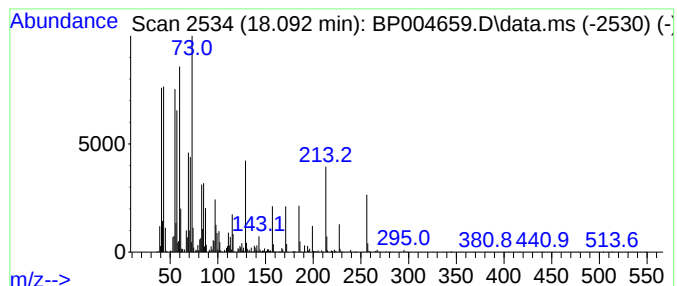
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	4.15 ng/ul	151782	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
3			Tridecanoic acid	214	C13H26O2	000638-53-9	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	91
5			Tridecanoic acid	214	C13H26O2	000638-53-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
Data File : BP004659.D
Acq On : 28 Jan 2021 15:30
Operator : CG/JU
Sample : M1203-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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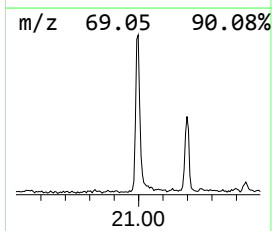
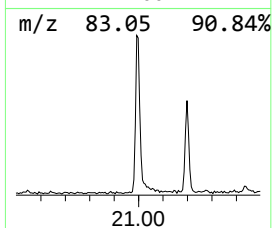
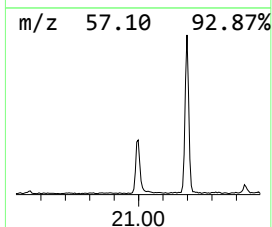
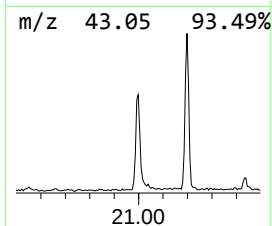
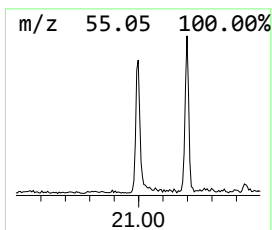
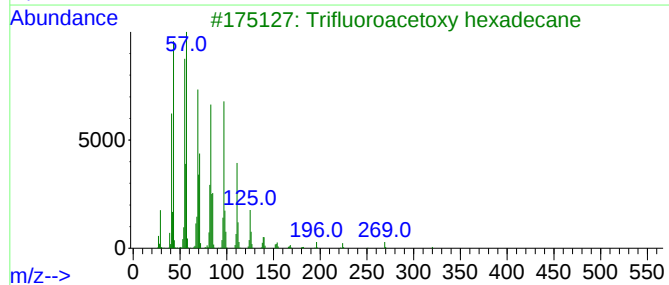
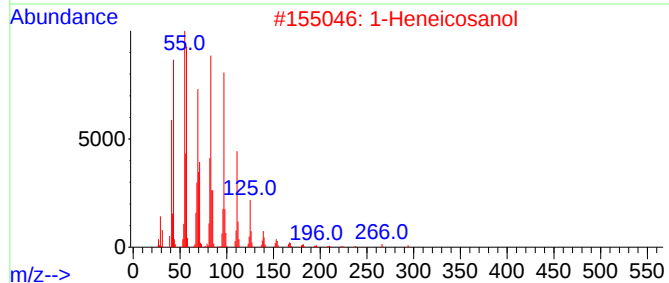
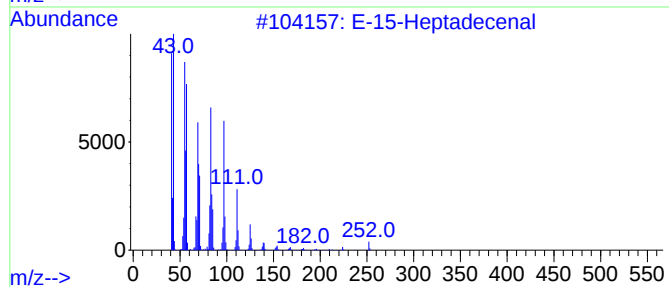
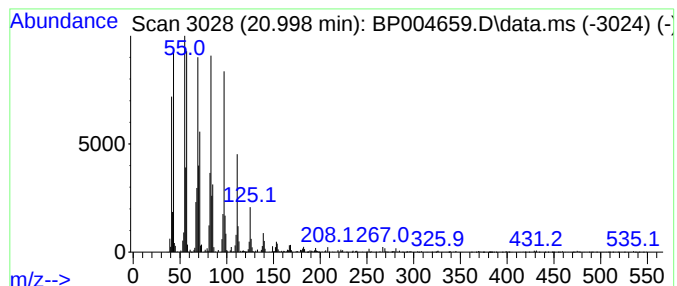
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 E-15-Heptadecenal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	2.65 ng/ul	319577	Chrysene-d12	21.280

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			E-15-Heptadecenal	252	C17H32O	1000130-97-9	96
2			1-Heneicosanol	312	C21H44O	015594-90-8	95
3			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
Data File : BP004659.D
Acq On : 28 Jan 2021 15:30
Operator : CG/JU
Sample : M1203-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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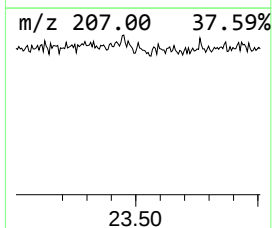
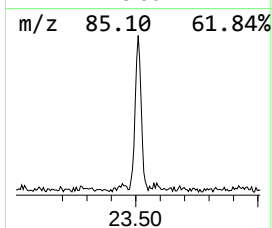
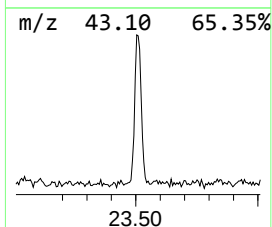
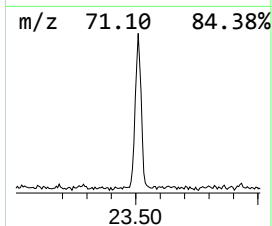
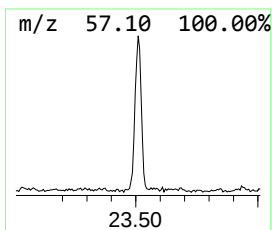
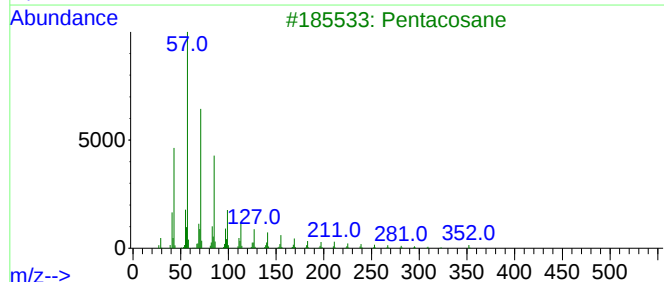
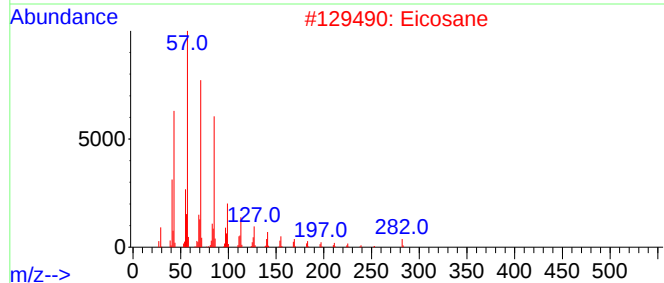
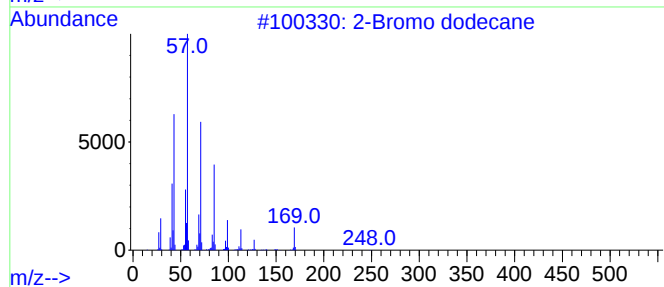
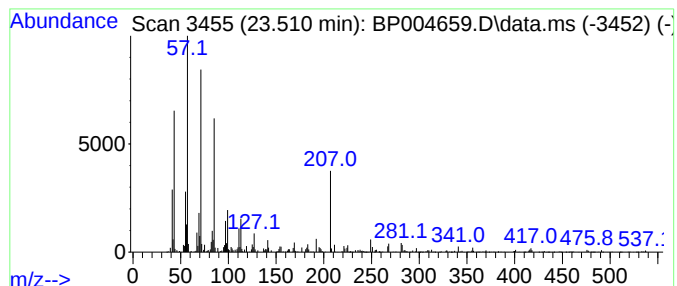
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Bromo dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.510	3.97 ng/ul	157382	Perylene-d12	23.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Bromo dodecane	248	C12H25Br	013187-99-0	81
2			Eicosane	282	C20H42	000112-95-8	81
3			Pentacosane	352	C25H52	000629-99-2	81
4			Eicosane	282	C20H42	000112-95-8	81
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
Data File : BP004659.D
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Sample : M1203-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

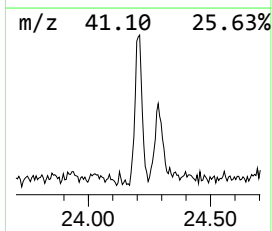
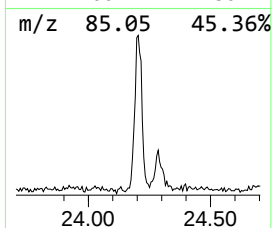
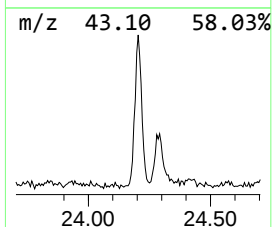
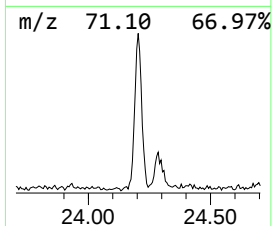
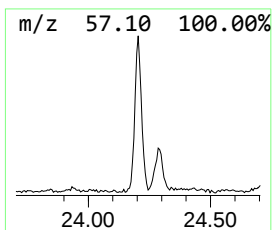
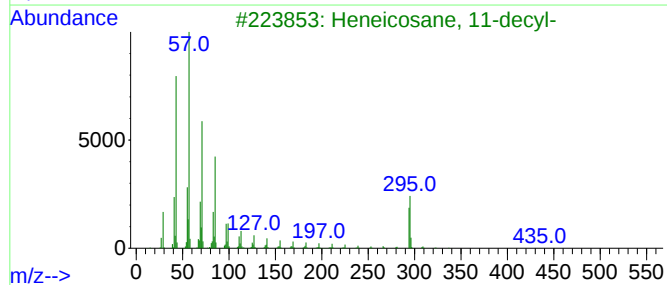
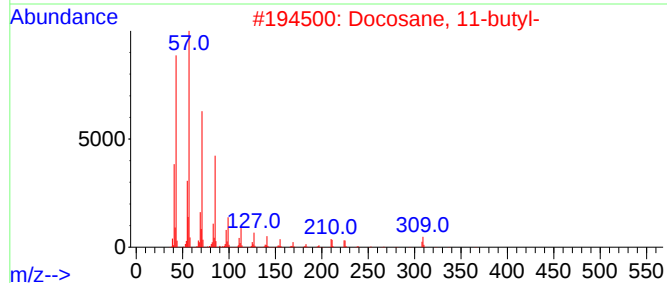
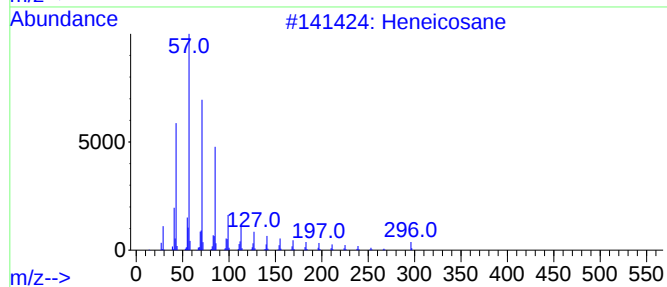
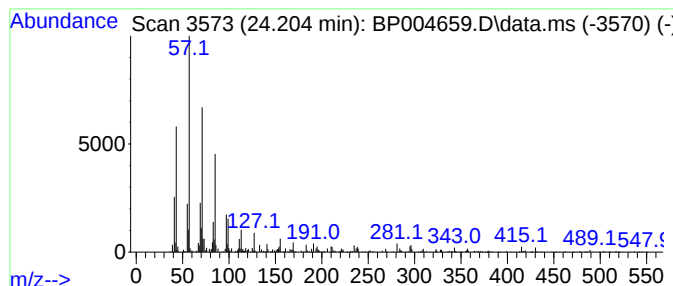
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP012221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.204	3.90 ng/ul	154871	Perylene-d12	23.598	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	93
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
4	Docosane, 5-butyl-	366	C26H54	055282-16-1	89
5	Tricosane	324	C23H48	000638-67-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012821\
Data File : BP004659.D
Acq On : 28 Jan 2021 15:30
Operator : CG/JU
Sample : M1203-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP012221.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	4.322	2.1	ng/ul	28104	1	7.793	267360	20.0
2-Pentanone, 4-...	4.916	9.2	ng/ul	122850	1	7.793	267360	20.0
o-Hydroxybiphenyl	14.716	25.7	ng/ul	703874	3	14.440	547321	20.0
n-Hexadecanoic ...	18.092	4.2	ng/ul	151782	4	17.192	731147	20.0
E-15-Heptadecenal	20.998	2.6	ng/ul	319577	5	21.280	2407480	20.0
2-Bromo dodecane	23.510	4.0	ng/ul	157382	6	23.598	793537	20.0
(DEL) Alkane: S...	24.204	3.9	ng/ul	154871	6	23.598	793537	20.0