

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017225.D
 Acq On : 19 Sep 2023 02:32
 Operator : MA/JU
 Sample : 04332-17
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AE4

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-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017225.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	35	38	44	rVB	32224	39940	1.37%	0.106%
2	3.940	140	145	150	rVB	74644	103497	3.55%	0.275%
3	4.058	162	165	169	rVB	12491	14427	0.49%	0.038%
4	4.446	227	231	236	rVB6	10171	15969	0.55%	0.042%
5	4.993	319	324	329	rBV	39875	57966	1.99%	0.154%
6	5.452	399	402	405	rVB5	4355	5015	0.17%	0.013%
7	5.581	418	424	429	rVB2	8843	14168	0.49%	0.038%
8	5.893	473	477	485	rVB2	5960	11895	0.41%	0.032%
9	6.093	508	511	513	rBV4	4729	4604	0.16%	0.012%
10	6.381	556	560	563	rVV6	3333	4803	0.16%	0.013%
11	6.440	568	570	573	rVB4	3448	3265	0.11%	0.009%
12	7.099	676	682	697	rBV	462156	794452	27.24%	2.109%
13	7.287	704	714	725	rVB	362047	547370	18.77%	1.453%
14	7.463	738	744	753	rBV	435858	704899	24.17%	1.871%
15	7.557	757	760	764	rVV6	5353	9184	0.31%	0.024%
16	7.616	766	770	773	rVB6	3316	5214	0.18%	0.014%
17	7.746	789	792	799	rVV8	4825	11349	0.39%	0.030%
18	7.946	819	826	837	rVB	752778	1175125	40.30%	3.120%
19	8.222	871	873	876	rVB4	3441	3412	0.12%	0.009%
20	8.434	905	909	912	rVB5	3478	4875	0.17%	0.013%
21	8.657	940	947	956	rBV	413085	677702	23.24%	1.799%
22	8.793	964	970	977	rBV	137989	223311	7.66%	0.593%
23	8.934	989	994	996	rBV5	5654	8751	0.30%	0.023%
24	8.999	1003	1005	1010	rVB6	2607	3337	0.11%	0.009%
25	9.146	1023	1030	1037	rBV	375616	632854	21.70%	1.680%
26	9.257	1047	1049	1054	rVB6	6689	8224	0.28%	0.022%
27	9.351	1061	1065	1072	rVB9	7422	13290	0.46%	0.035%
28	9.857	1143	1151	1160	rBV	303629	496975	17.04%	1.319%
29	9.998	1169	1175	1179	rBV2	27066	49665	1.70%	0.132%
30	10.116	1193	1195	1202	rVB8	4555	7998	0.27%	0.021%
31	10.222	1208	1213	1219	rBV3	12791	21702	0.74%	0.058%
32	10.387	1234	1241	1251	rBV	433210	735281	25.21%	1.952%
33	10.510	1259	1262	1266	rBV6	5447	7882	0.27%	0.021%
34	10.769	1298	1306	1314	rBV	922051	1539660	52.80%	4.088%
35	10.934	1326	1334	1345	rBV	195952	368127	12.62%	0.977%
36	11.334	1394	1402	1408	rBV	6560	16119	0.55%	0.043%

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Title : SVOA CALIBRATION

37	11.516	1429	1433	1434	rBV4	4213	3782	0.13%	0.010%
38	11.593	1440	1446	1455	rVB3	10033	25039	0.86%	0.066%
39	11.651	1455	1456	1460	rBV4	3031	3262	0.11%	0.009%
40	11.804	1479	1482	1487	rVB7	2359	3639	0.12%	0.010%
41	12.122	1532	1536	1539	rBV6	2574	3809	0.13%	0.010%
42	12.240	1553	1556	1557	rBV3	4607	3798	0.13%	0.010%
43	12.351	1568	1575	1579	rBV6	9470	17121	0.59%	0.045%
44	12.522	1598	1604	1608	rBV7	11794	22762	0.78%	0.060%
45	12.645	1622	1625	1630	rVB7	3097	5599	0.19%	0.015%
46	12.716	1634	1637	1640	rBV5	2458	3758	0.13%	0.010%
47	12.816	1650	1654	1656	rBV5	2940	3590	0.12%	0.010%
48	12.863	1656	1662	1663	rBV5	2902	5031	0.17%	0.013%
49	13.045	1691	1693	1698	rVB6	2327	3165	0.11%	0.008%
50	13.357	1741	1746	1749	rBV5	8018	9095	0.31%	0.024%
51	13.445	1754	1761	1768	rVB2	36370	60162	2.06%	0.160%
52	13.539	1772	1777	1778	rBV5	3358	4922	0.17%	0.013%
53	13.639	1787	1794	1802	rBV	87652	145143	4.98%	0.385%
54	13.792	1813	1820	1825	rBV	7208	15920	0.55%	0.042%
55	13.910	1838	1840	1846	rVB7	4061	6976	0.24%	0.019%
56	14.016	1852	1858	1870	rBV	697232	944815	32.40%	2.508%
57	14.304	1901	1907	1921	rBV2	838459	1232360	42.26%	3.272%
58	14.434	1926	1929	1933	rVV6	5252	5872	0.20%	0.016%
59	14.604	1950	1958	1972	rBV	1273778	1816200	62.28%	4.822%
60	14.834	1991	1997	2011	rBV2	110142	222710	7.64%	0.591%
61	15.004	2020	2026	2030	rBV8	9039	17275	0.59%	0.046%
62	15.039	2030	2032	2038	rVB7	5455	8206	0.28%	0.022%
63	15.298	2073	2076	2080	rBV6	4650	7298	0.25%	0.019%
64	15.345	2082	2084	2087	rVB4	4214	4235	0.15%	0.011%
65	15.386	2089	2091	2095	rVV5	4037	4170	0.14%	0.011%
66	15.457	2102	2103	2106	rBV3	3351	3466	0.12%	0.009%
67	15.598	2119	2127	2144	rBV	1055216	1527677	52.38%	4.056%
68	15.733	2144	2150	2157	rBV	156344	227344	7.80%	0.604%
69	15.833	2165	2167	2168	rBV2	5228	4651	0.16%	0.012%
70	15.951	2185	2187	2190	rBV4	4382	4665	0.16%	0.012%
71	16.263	2236	2240	2244	rVB7	7197	10012	0.34%	0.027%
72	16.304	2244	2247	2253	rVB8	4691	8371	0.29%	0.022%
73	16.392	2257	2262	2264	rBV3	12300	17047	0.58%	0.045%
74	16.428	2264	2268	2274	rVB6	20087	31535	1.08%	0.084%
75	16.692	2311	2313	2315	rBV3	3983	3827	0.13%	0.010%
76	16.722	2315	2318	2323	rVV5	15480	22577	0.77%	0.060%

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 Title : SVOA CALIBRATION

77	16.833	2335	2337	2341	rVB5	4245	5331	0.18%	0.014%
78	17.227	2399	2404	2410	rBV	106469	173084	5.94%	0.460%
79	17.292	2411	2415	2422	rVV3	88194	154825	5.31%	0.411%
80	17.375	2423	2429	2440	rVV	1569295	2237082	76.71%	5.939%
81	17.475	2440	2446	2455	rVV2	1182561	1701671	58.35%	4.518%
82	18.110	2548	2554	2560	rVV4	114793	201146	6.90%	0.534%
83	18.163	2560	2563	2568	rVB	125562	152761	5.24%	0.406%
84	18.216	2568	2572	2586	rVB	645711	915453	31.39%	2.431%
85	18.504	2617	2621	2629	rBV7	79056	137859	4.73%	0.366%
86	19.039	2708	2712	2715	rBV	141876	175494	6.02%	0.466%
87	19.345	2761	2764	2765	rBV3	76285	82232	2.82%	0.218%
88	19.457	2779	2783	2787	rBV	185213	267406	9.17%	0.710%
89	19.539	2794	2797	2806	rVB	470547	653009	22.39%	1.734%
90	19.716	2822	2827	2835	rVB	1624002	2048296	70.24%	5.438%
91	20.180	2903	2906	2908	rBV2	85857	103753	3.56%	0.275%
92	20.980	3039	3042	3046	rVB2	161931	192649	6.61%	0.511%
93	21.151	3068	3071	3078	rVB2	303662	475802	16.32%	1.263%
94	21.480	3122	3127	3132	rVB	1916831	2506303	85.94%	6.654%
95	22.092	3228	3231	3241	rVB2	360428	671229	23.02%	1.782%
96	23.133	3403	3408	3416	rVB	1029496	1696319	58.17%	4.504%
97	23.227	3419	3424	3441	rVV2	605901	1789510	61.36%	4.751%
98	23.833	3522	3527	3534	rVB2	1091285	2081931	71.39%	5.527%
99	24.004	3550	3556	3568	rVB	1384885	2916296	100.00%	7.743%
100	24.557	3645	3650	3661	rVB	707997	1512387	51.86%	4.015%

Sum of corrected areas: 37665021

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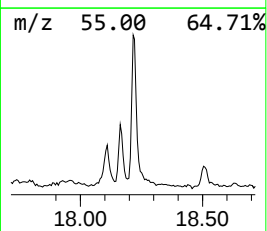
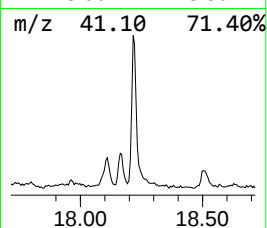
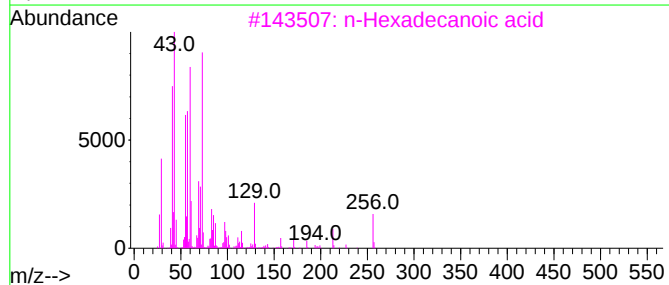
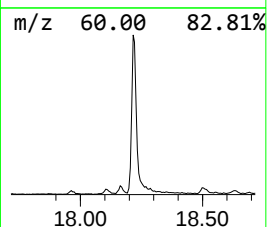
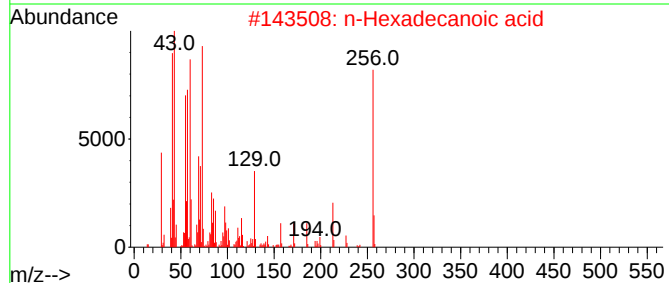
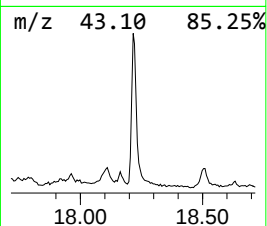
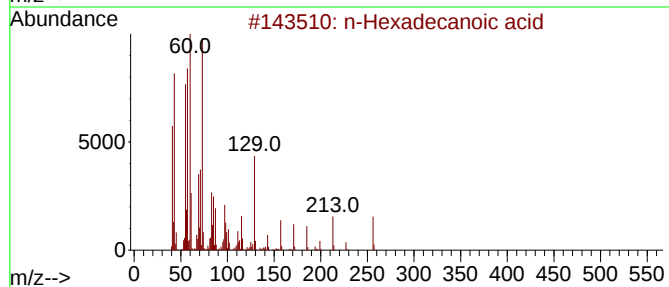
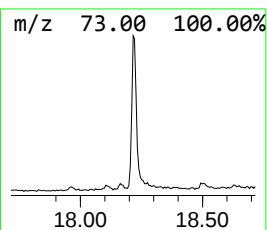
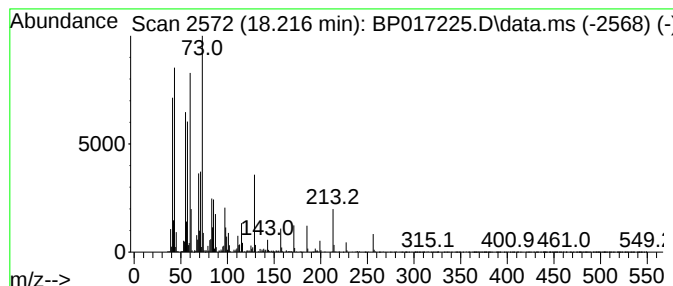
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	8.18 ng/ul	915453	Phenanthrene-d10	17.375

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



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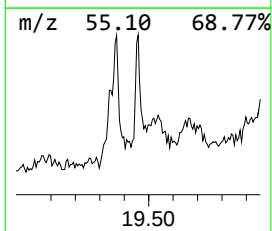
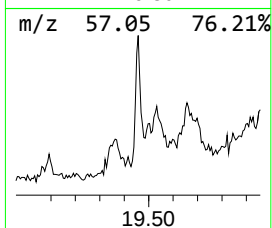
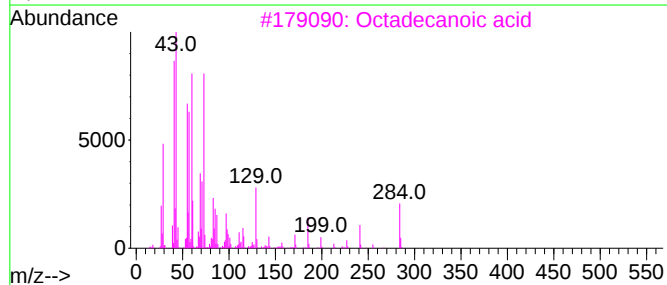
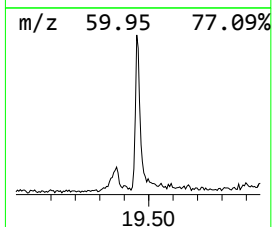
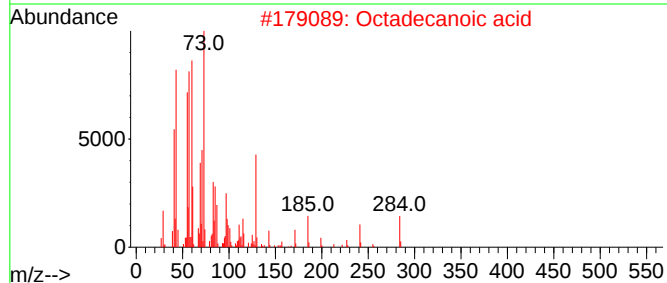
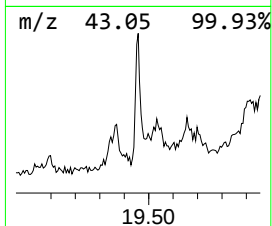
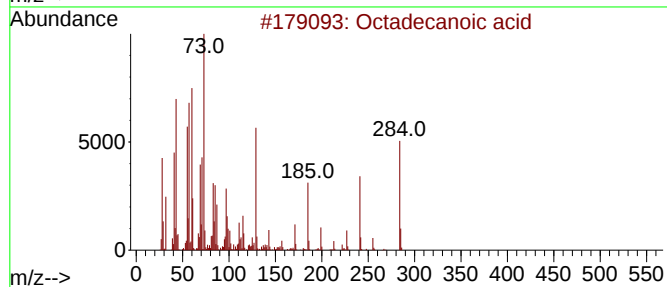
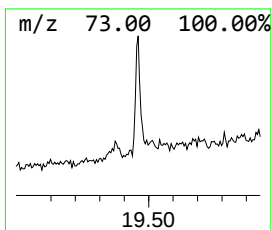
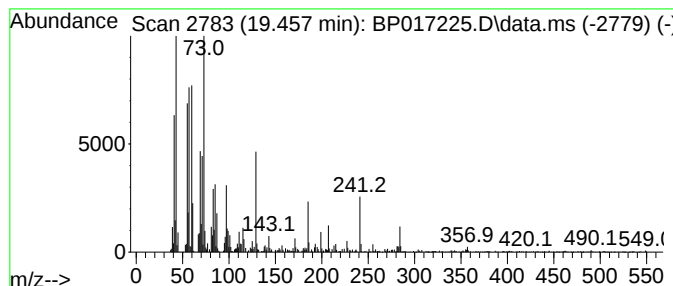
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	2.13 ng/ul	267406	Chrysene-d12	21.480

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		Octadecanoic acid	284	C18H36O2	000057-11-4	97
4		Octadecanoic acid	284	C18H36O2	000057-11-4	95
5		Octadecanoic acid	284	C18H36O2	000057-11-4	93



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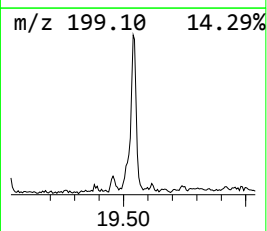
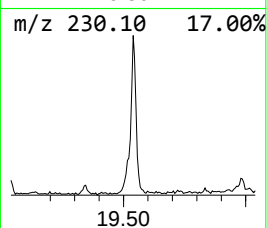
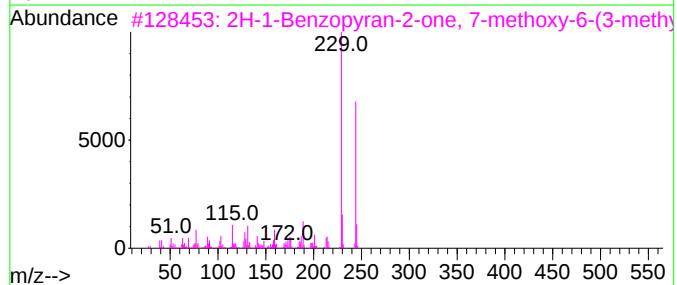
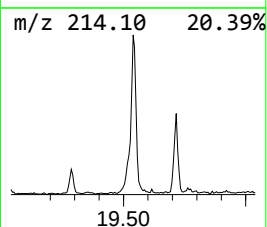
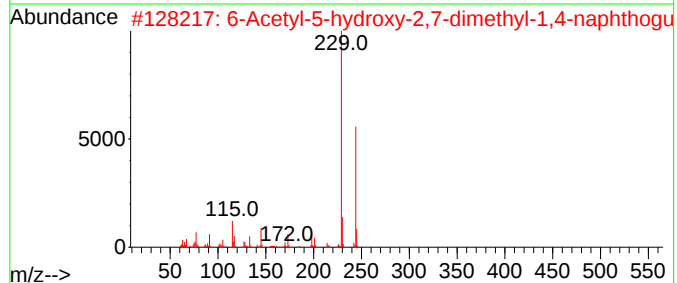
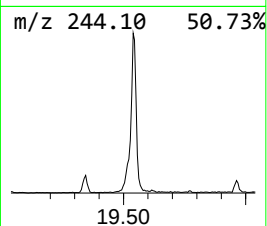
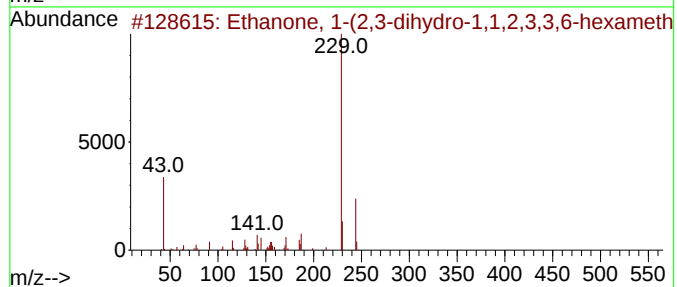
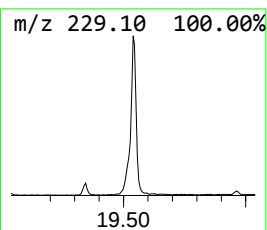
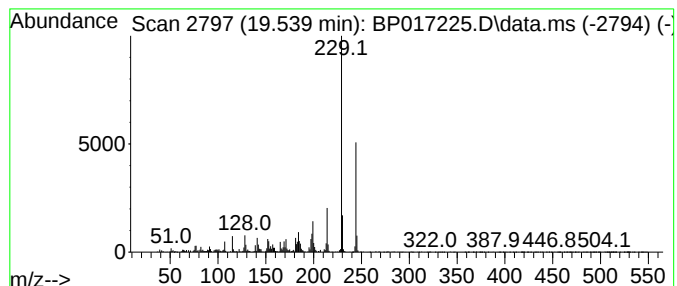
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanone, 1-(2,3-dihydro-1,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	5.21 ng/ul	653009	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-(2,3-dihydro-1,1,2,3...	244	C17H24O	015323-35-0	70
2			6-Acetyl-5-hydroxy-2,7-dimethyl-...	244	C14H12O4	001086-07-3	68
3			2H-1-Benzopyran-2-one, 7-methoxy...	244	C15H16O3	000581-31-7	64
4			Cyclopentanone, tetrakis(1-methy...	244	C17H24O	096446-08-1	64
5			2-Propenoic acid, 2-cyano-3-[4-(...	244	C14H16N2O2	1000115-72-8	62



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ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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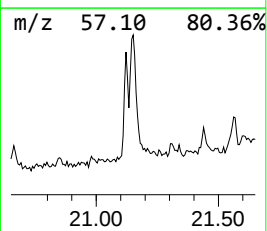
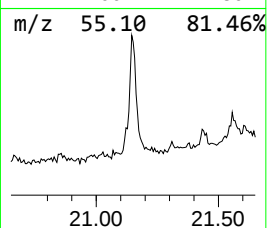
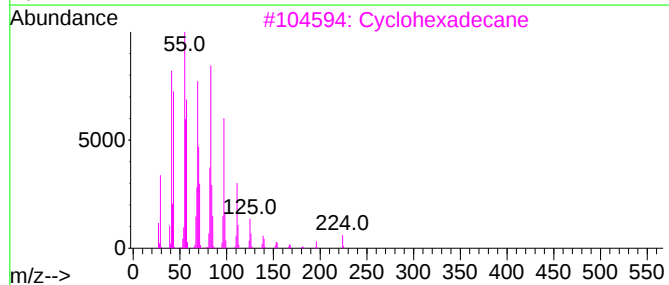
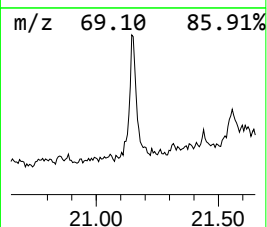
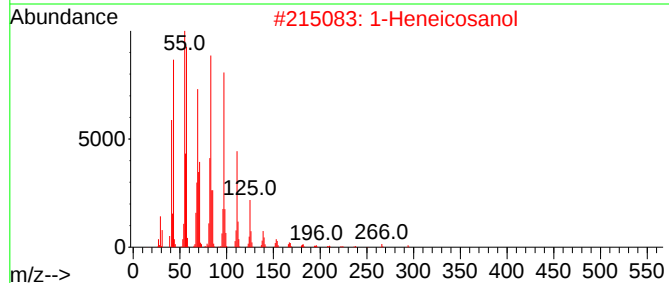
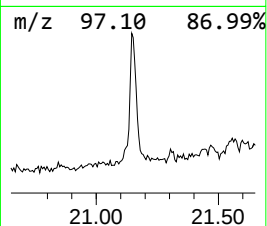
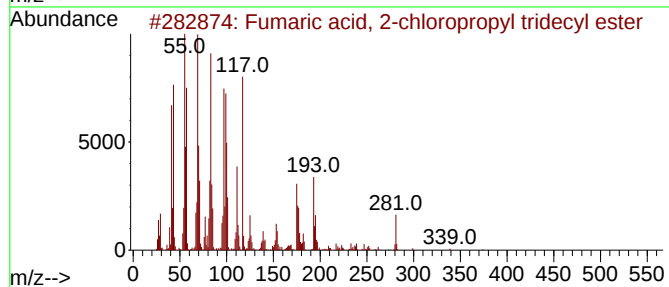
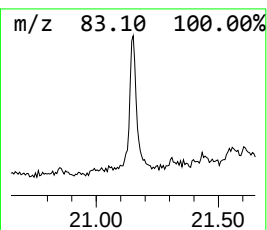
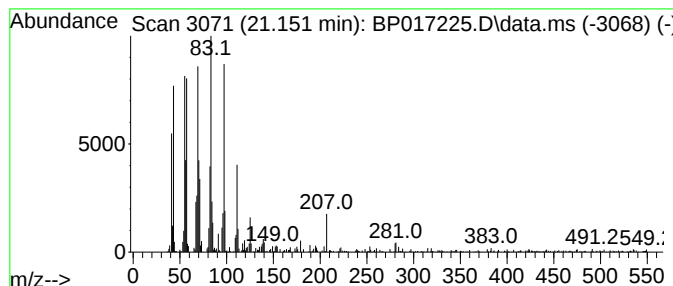
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Fumaric acid, 2-chloropropyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.151	3.80 ng/ul	475802	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	96
2			1-Heneicosanol	312	C21H44O	015594-90-8	94
3			Cyclohexadecane	224	C16H32	000295-65-8	93
4			Behenic alcohol	326	C22H46O	000661-19-8	91
5			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017225.D
Acq On : 19 Sep 2023 02:32
Operator : MA/JU
Sample : 04332-17
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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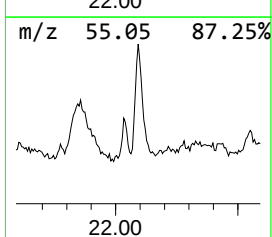
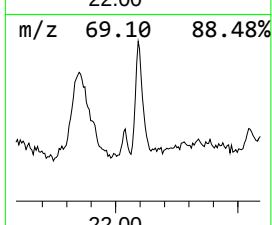
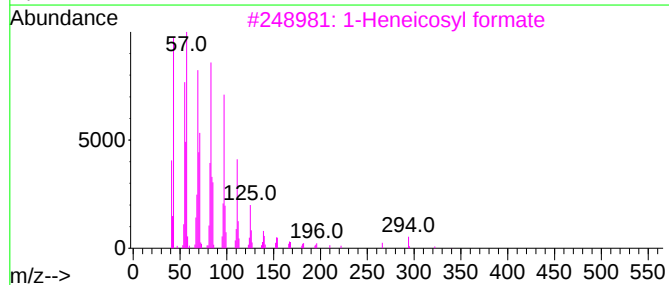
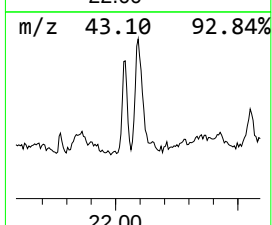
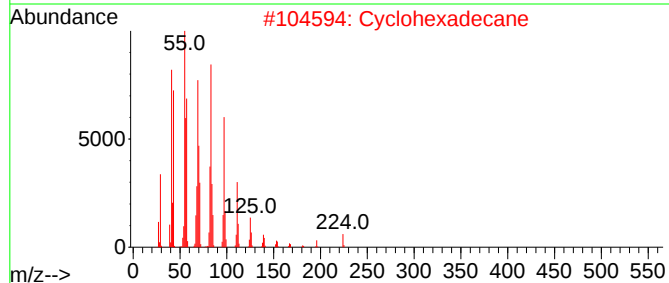
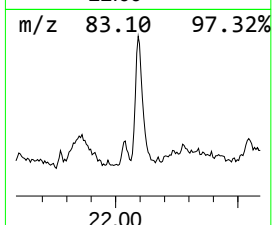
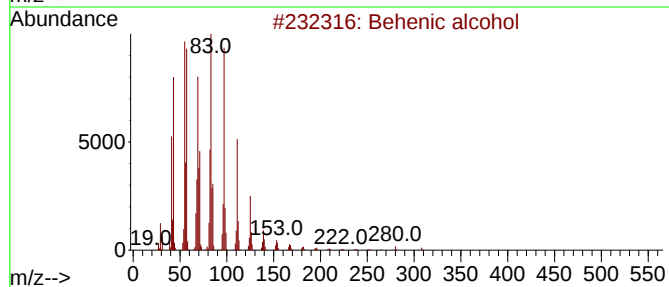
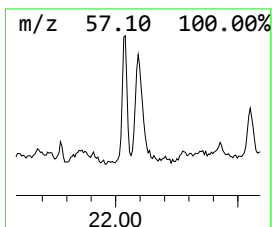
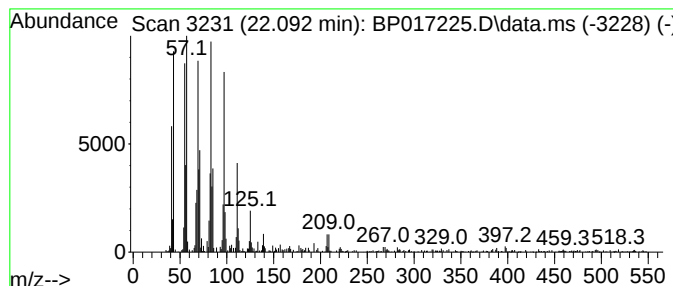
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Behenic alcohol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.092	5.36 ng/ul	671229	Chrysene-d12	21.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	93
2			Cyclohexadecane	224	C16H32	000295-65-8	93
3			1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5			1-Heptacosanol	396	C27H56O	002004-39-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017225.D
Acq On : 19 Sep 2023 02:32
Operator : MA/JU
Sample : 04332-17
Misc :
ALS Vial : 29 Sample Multiplier: 1

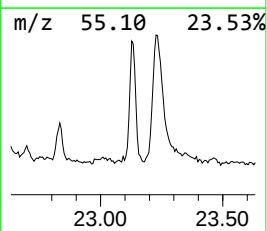
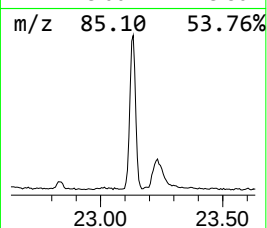
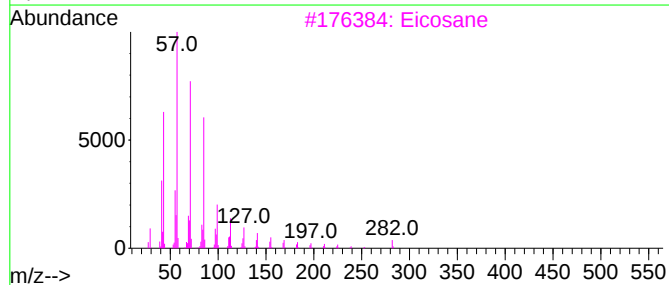
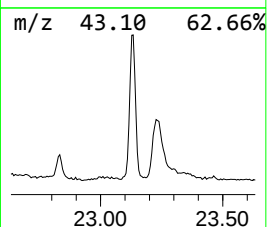
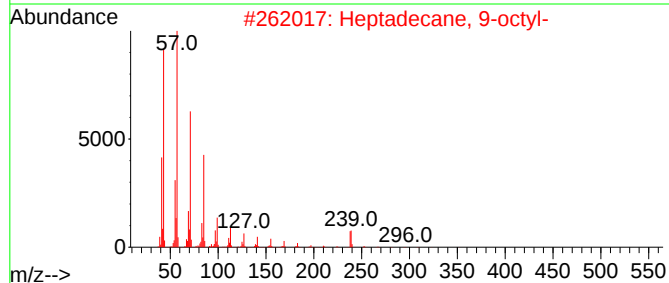
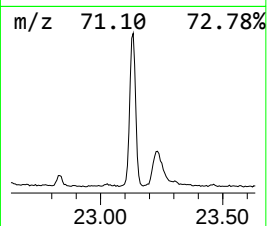
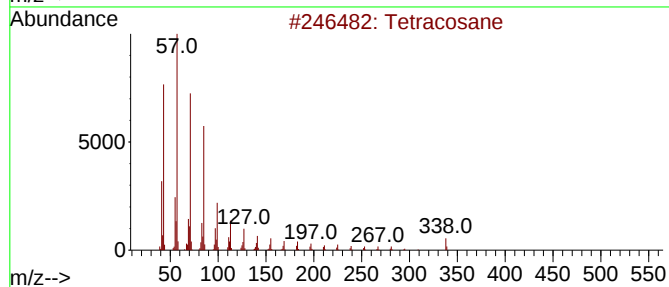
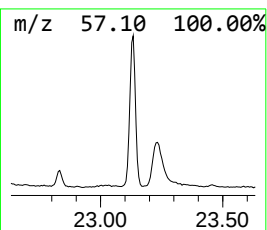
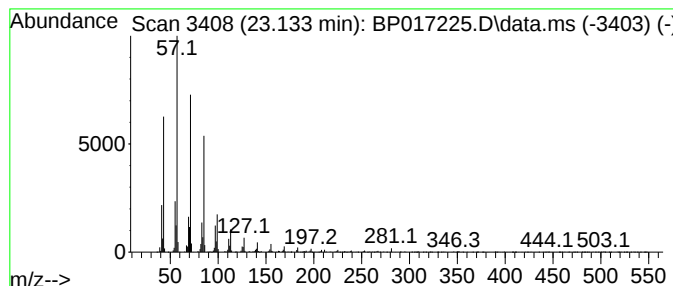
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.133	11.63 ng/ul	1696320	Perylene-d12	24.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3	Eicosane	282	C20H42	000112-95-8	94
4	Heptadecane	240	C17H36	000629-78-7	93
5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017225.D
Acq On : 19 Sep 2023 02:32
Operator : MA/JU
Sample : 04332-17
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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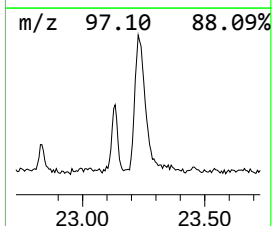
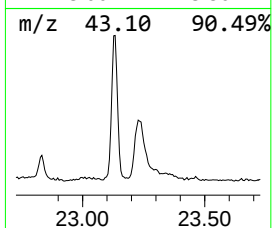
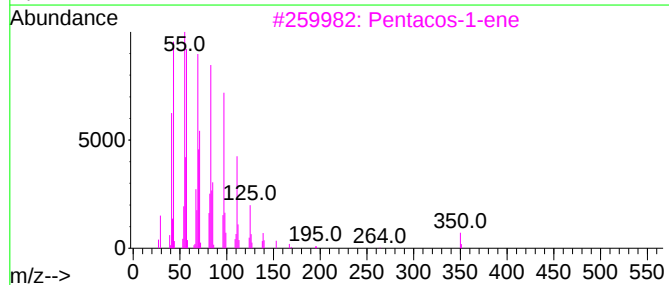
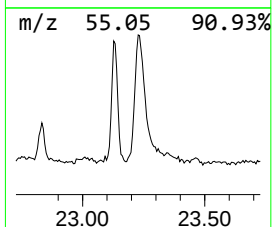
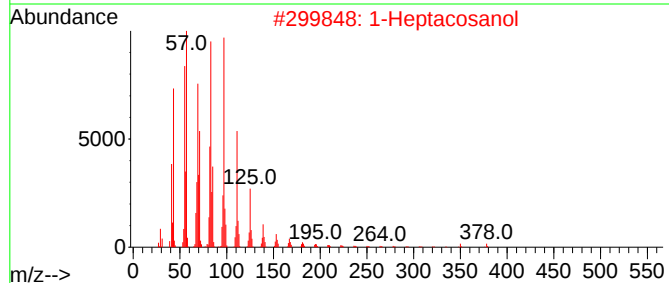
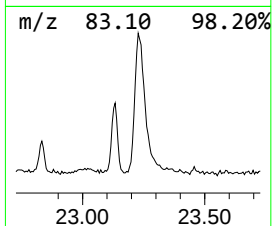
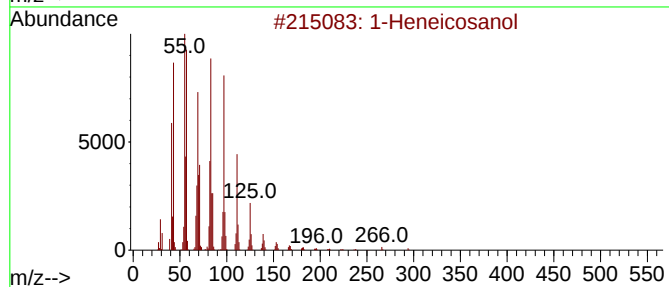
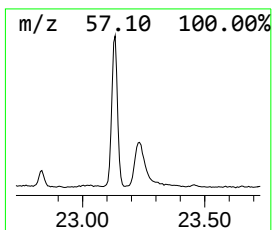
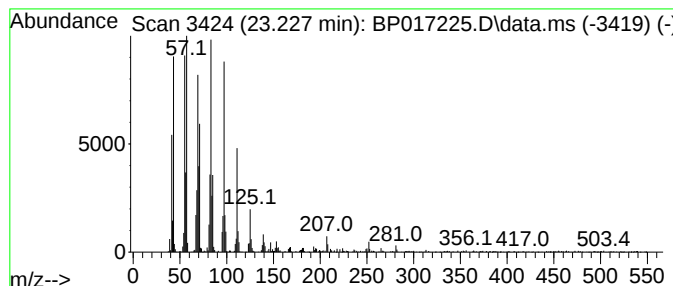
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	12.27 ng/ul	1789510	Perylene-d12	24.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol			312	C21H44O	015594-90-8	95
2	1-Heptacosanol			396	C27H56O	002004-39-9	94
3	Pentacos-1-ene			350	C25H50	016980-85-1	94
4	Nonacos-1-ene			406	C29H58	018835-35-3	94
5	Heptacos-1-ene			378	C27H54	015306-27-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017225.D
Acq On : 19 Sep 2023 02:32
Operator : MA/JU
Sample : 04332-17
Misc :
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ClientSampleId :
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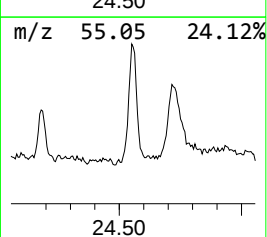
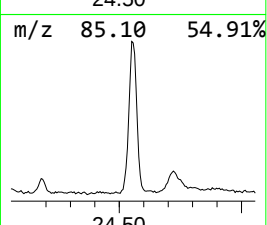
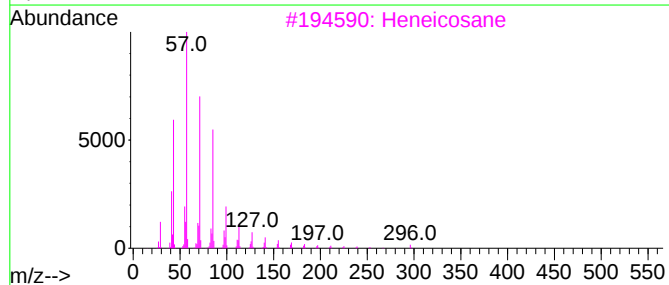
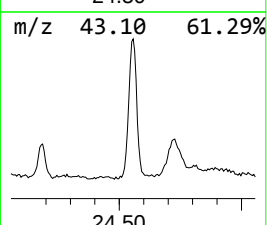
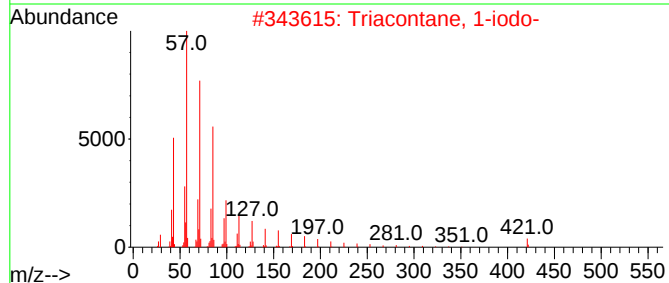
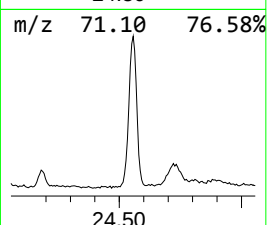
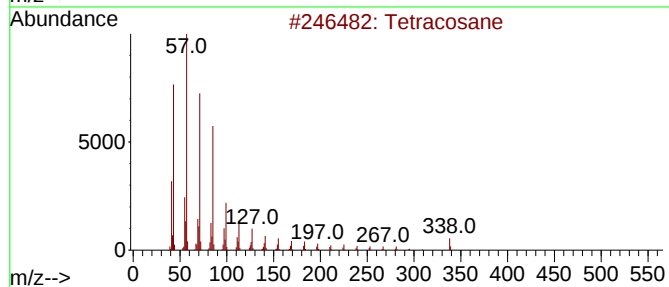
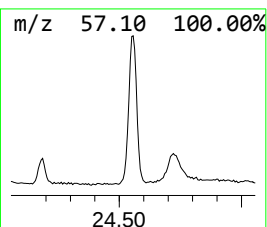
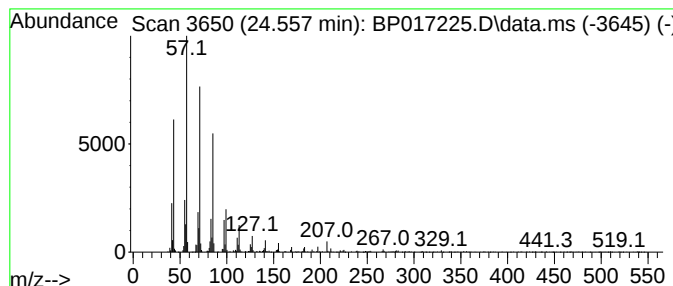
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.557	10.37 ng/ul	1512390	Perylene-d12	24.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	94
2			Triacotane, 1-iodo-	548	C30H61I	1000406-32-3	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Triacotane	422	C30H62	000638-68-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
Data File : BP017225.D
Acq On : 19 Sep 2023 02:32
Operator : MA/JU
Sample : 04332-17
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.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
n-Hexadecanoic ...	18.216	8.2	ng/ul	915453	4	17.375	2237080	20.0
Octadecanoic acid	19.457	2.1	ng/ul	267406	5	21.480	2506300	20.0
Ethanone, 1-(2,...	19.539	5.2	ng/ul	653009	5	21.480	2506300	20.0
Fumaric acid, 2...	21.151	3.8	ng/ul	475802	5	21.480	2506300	20.0
Behenic alcohol	22.092	5.4	ng/ul	671229	5	21.480	2506300	20.0
(DEL) Alkane: S...	23.133	11.6	ng/ul	1696320	6	24.004	2916300	20.0
1-Heneicosanol	23.227	12.3	ng/ul	1789510	6	24.004	2916300	20.0
(DEL) Alkane: S...	24.557	10.4	ng/ul	1512390	6	24.004	2916300	20.0