

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
 Data File : BP020747.D
 Acq On : 02 Jul 2024 12:19
 Operator : MA/JU
 Sample : P2963-02
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4B47

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-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020747.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	44	48	56	rVB	56292	79642	2.05%	0.114%
2	3.546	91	95	104	rVB	38308	57827	1.49%	0.083%
3	3.693	117	120	130	rVV	90786	134435	3.46%	0.193%
4	3.787	131	136	140	rVB	44968	61027	1.57%	0.088%
5	3.993	166	171	179	rVB	343768	502314	12.93%	0.720%
6	4.546	260	265	268	rVV	39620	56394	1.45%	0.081%
7	4.587	268	272	279	rVB	40163	54689	1.41%	0.078%
8	4.893	317	324	332	rBV	200376	287056	7.39%	0.412%
9	5.852	480	487	499	rBV	31659	63844	1.64%	0.092%
10	6.758	632	641	650	rBV6	12303	38418	0.99%	0.055%
11	6.922	660	669	683	rVB	1002256	1626721	41.89%	2.333%
12	7.081	690	696	710	rBV	776457	1229361	31.65%	1.763%
13	7.281	718	730	738	rBV	1062339	1639253	42.21%	2.351%
14	7.381	744	747	752	rVB	41126	61968	1.60%	0.089%
15	7.734	797	807	814	rBV	891303	1326594	34.16%	1.902%
16	7.805	814	819	823	rVV2	37637	70400	1.81%	0.101%
17	8.099	864	869	879	rBV	18953	37712	0.97%	0.054%
18	8.434	920	926	934	rBV	1343386	2004915	51.62%	2.875%
19	8.552	942	946	956	rVB	22505	37555	0.97%	0.054%
20	8.887	996	1003	1016	rVV	958990	1573471	40.52%	2.256%
21	9.263	1060	1067	1075	rVB	29744	50167	1.29%	0.072%
22	9.446	1093	1098	1107	rVB	124356	192343	4.95%	0.276%
23	9.593	1116	1123	1137	rBV	740059	1357824	34.96%	1.947%
24	10.140	1208	1216	1225	rBV	1220005	2052488	52.85%	2.943%
25	10.204	1225	1227	1235	rVV3	18070	39252	1.01%	0.056%
26	10.510	1270	1279	1284	rVV	1079229	1867464	48.09%	2.678%
27	10.557	1284	1287	1294	rVB	160701	254470	6.55%	0.365%
28	10.646	1295	1302	1320	rBV	670242	1203204	30.98%	1.725%
29	11.046	1363	1370	1379	rBV6	34939	73579	1.89%	0.106%
30	11.252	1399	1405	1418	rBV	112248	249292	6.42%	0.357%
31	11.899	1508	1515	1523	rBV3	60689	151652	3.90%	0.217%
32	12.093	1539	1548	1555	rVV2	23953	61234	1.58%	0.088%
33	12.169	1555	1561	1566	rVB	25567	47291	1.22%	0.068%
34	12.387	1592	1598	1603	rVV	28281	44066	1.13%	0.063%
35	13.122	1716	1723	1727	rBV10	13356	33456	0.86%	0.048%
36	13.175	1727	1732	1739	rVV4	32952	64162	1.65%	0.092%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	13.545	1786	1795	1801	rBV9	18166	45119	1.16%	0.065%
38	13.681	1813	1818	1823	rVB2	31204	53154	1.37%	0.076%
39	13.781	1828	1835	1843	rVV	1908566	2808755	72.32%	4.028%
40	14.057	1876	1882	1900	rVB	2108754	3424547	88.18%	4.911%
41	14.369	1926	1935	1941	rBV	1504574	2333562	60.09%	3.346%
42	14.434	1941	1946	1952	rVB	91654	125368	3.23%	0.180%
43	14.493	1952	1956	1963	rVB3	48373	82193	2.12%	0.118%
44	14.593	1967	1973	1984	rBV2	586187	1124004	28.94%	1.612%
45	14.740	1994	1998	2000	rBV3	25551	36077	0.93%	0.052%
46	14.775	2000	2004	2015	rVB	66442	143123	3.69%	0.205%
47	15.169	2067	2071	2077	rBV7	21821	44346	1.14%	0.064%
48	15.381	2101	2107	2112	rBV	2730622	3802399	97.91%	5.452%
49	15.434	2112	2116	2121	rVB	119354	168333	4.33%	0.241%
50	15.504	2123	2128	2136	rBV	653202	906464	23.34%	1.300%
51	15.645	2147	2152	2158	rVB	290119	441575	11.37%	0.633%
52	15.740	2164	2168	2176	rVB	35266	61716	1.59%	0.088%
53	15.887	2190	2193	2201	rVB2	38498	76048	1.96%	0.109%
54	15.975	2201	2208	2214	rBV9	17007	49451	1.27%	0.071%
55	16.134	2228	2235	2239	rVB5	37371	83862	2.16%	0.120%
56	16.269	2253	2258	2263	rBV2	64157	101218	2.61%	0.145%
57	16.398	2274	2280	2285	rBV	285180	409815	10.55%	0.588%
58	16.457	2286	2290	2296	rVV7	42747	84775	2.18%	0.122%
59	16.704	2327	2332	2337	rBV2	89748	140341	3.61%	0.201%
60	16.834	2349	2354	2360	rVB2	43365	71958	1.85%	0.103%
61	16.934	2368	2371	2374	rVB3	28548	34032	0.88%	0.049%
62	16.986	2377	2380	2387	rVB	88436	129507	3.33%	0.186%
63	17.057	2388	2392	2395	rBV2	102442	132633	3.42%	0.190%
64	17.092	2395	2398	2404	rVB	164760	205866	5.30%	0.295%
65	17.186	2406	2414	2417	rBV	1575804	2412741	62.13%	3.460%
66	17.228	2417	2421	2425	rVB	1177458	1557419	40.10%	2.233%
67	17.281	2425	2430	2434	rBV	2610940	3440241	88.58%	4.933%
68	17.363	2440	2444	2453	rVB2	144105	219394	5.65%	0.315%
69	17.610	2482	2486	2489	rBV	50910	70702	1.82%	0.101%
70	17.651	2490	2493	2496	rVV	59584	68209	1.76%	0.098%
71	17.792	2512	2517	2525	rVB2	185202	399008	10.27%	0.572%
72	17.922	2536	2539	2546	rVB3	69061	119061	3.07%	0.171%
73	18.116	2564	2572	2574	rBV2	267066	573616	14.77%	0.823%
74	18.222	2587	2590	2596	rVB2	44033	60447	1.56%	0.087%
75	18.292	2596	2602	2606	rBV2	292522	501672	12.92%	0.719%
76	18.398	2616	2620	2626	rBV3	54452	98642	2.54%	0.141%

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 Title : SVOA CALIBRATION

77	18.610	2649	2656	2660	rBV	102676	209053	5.38%	0.300%
78	18.657	2660	2664	2669	rVB2	127195	185387	4.77%	0.266%
79	18.863	2696	2699	2703	rVB	79669	94681	2.44%	0.136%
80	18.916	2705	2708	2711	rVV	46326	63167	1.63%	0.091%
81	19.045	2726	2730	2734	rBV	101408	183923	4.74%	0.264%
82	19.092	2734	2738	2744	rVV5	88992	199489	5.14%	0.286%
83	19.186	2751	2754	2760	rVB3	77588	108568	2.80%	0.156%
84	19.316	2771	2776	2783	rVB	1603993	2214832	57.03%	3.176%
85	19.457	2797	2800	2807	rBV5	126182	224766	5.79%	0.322%
86	19.669	2830	2836	2839	rBV	2399769	3883637	100.00%	5.569%
87	19.698	2839	2841	2850	rVB	1122017	1333114	34.33%	1.912%
88	20.222	2926	2930	2934	rVB	207924	306072	7.88%	0.439%
89	20.257	2934	2936	2939	rBV3	94596	111911	2.88%	0.160%
90	20.322	2944	2947	2951	rBV	178631	222977	5.74%	0.320%
91	20.380	2954	2957	2961	rVB2	222234	291126	7.50%	0.417%
92	21.316	3111	3116	3122	rBV3	998515	1515759	39.03%	2.173%
93	21.557	3153	3157	3161	rVV	743263	963621	24.81%	1.382%
94	21.633	3163	3170	3176	rVV3	1372353	3424924	88.19%	4.911%
95	21.686	3176	3179	3191	rVB	593104	1141974	29.40%	1.637%
96	22.533	3319	3323	3325	rBV2	455648	727611	18.74%	1.043%
97	23.933	3556	3561	3569	rBV	414241	1144945	29.48%	1.642%
98	24.133	3590	3595	3602	rVB	598174	1241886	31.98%	1.781%
99	24.768	3696	3703	3710	rBV2	839128	2108834	54.30%	3.024%
100	24.998	3735	3742	3752	rVB	992180	2509987	64.63%	3.599%

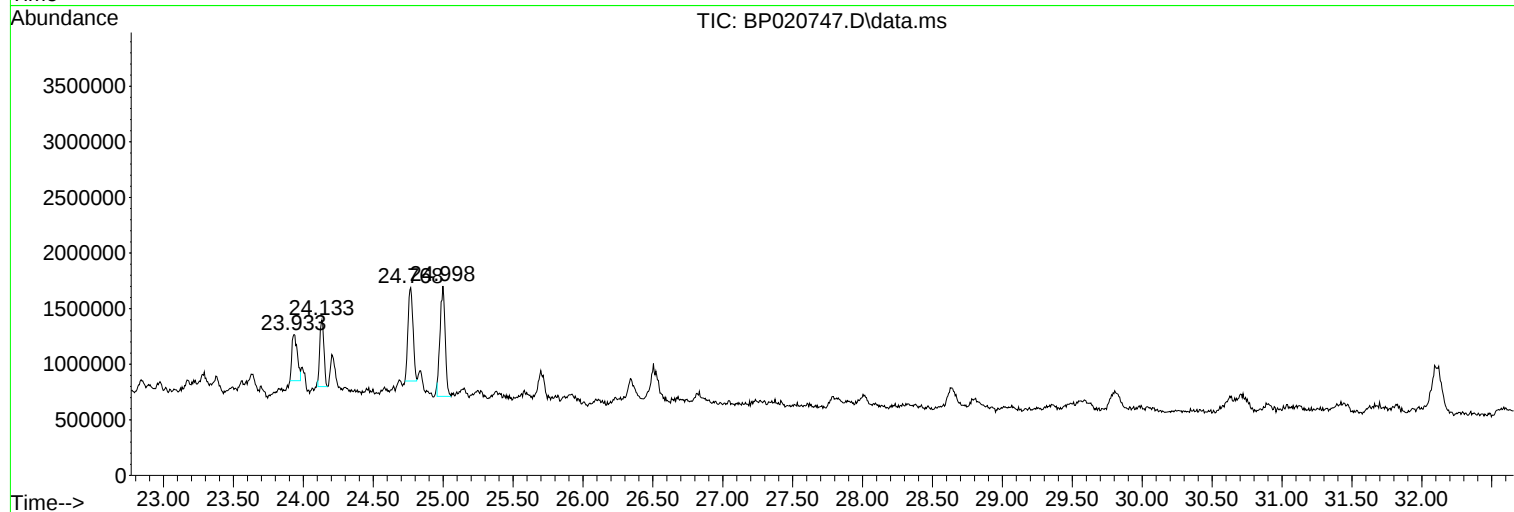
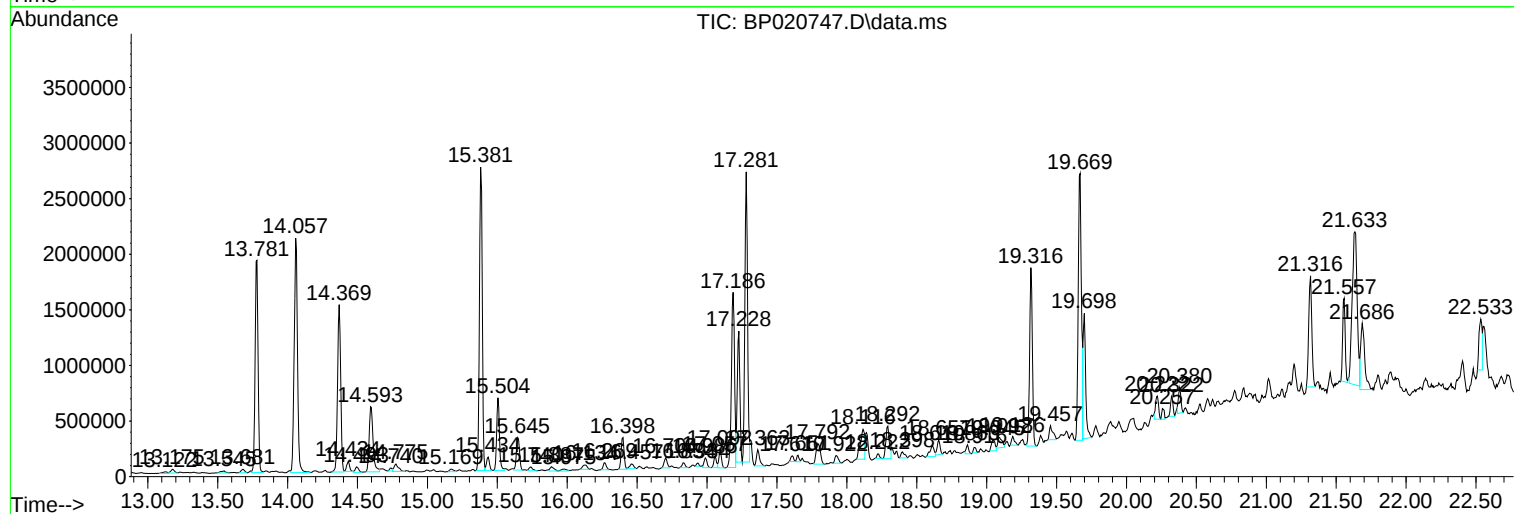
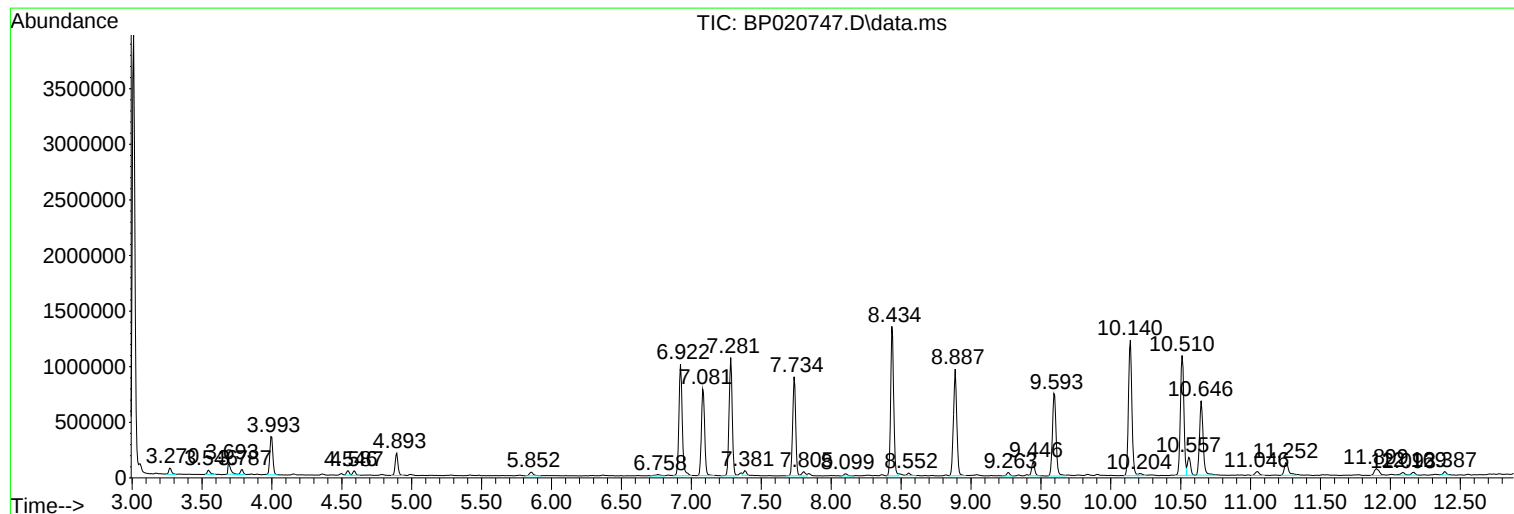
Sum of corrected areas: 69739177

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Misc :
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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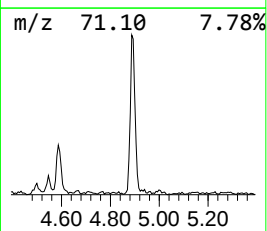
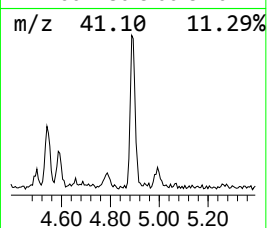
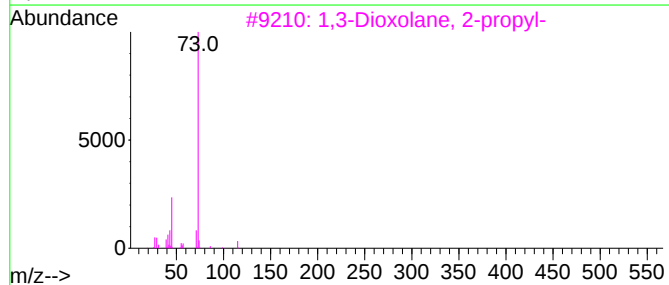
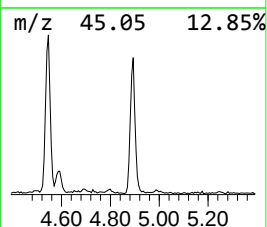
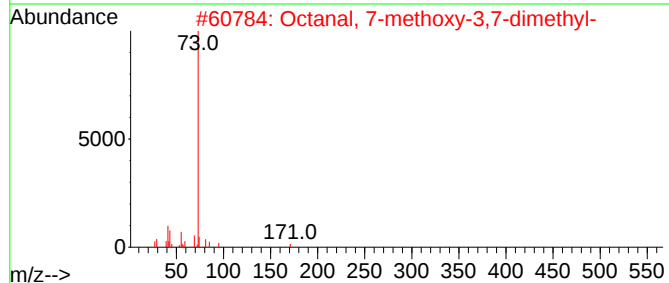
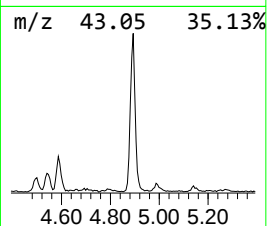
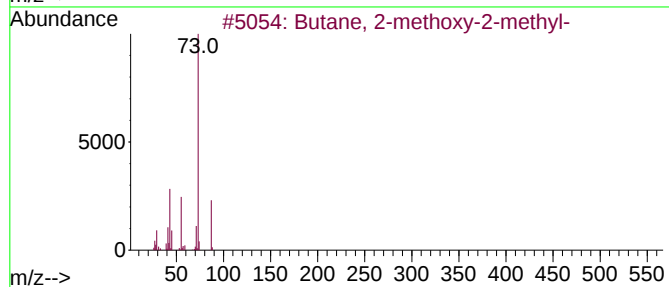
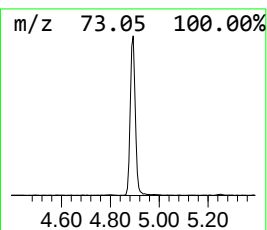
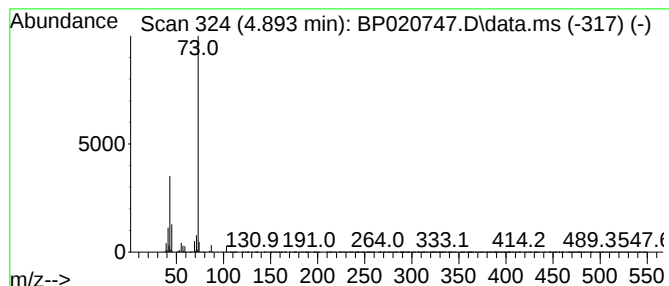
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.893	4.33 ng/ul	287056	1,4-Dichlorobenzene-d4	7.734

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	33
2			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	28
3			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9



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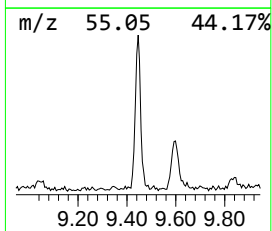
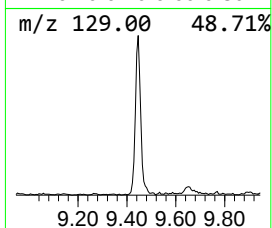
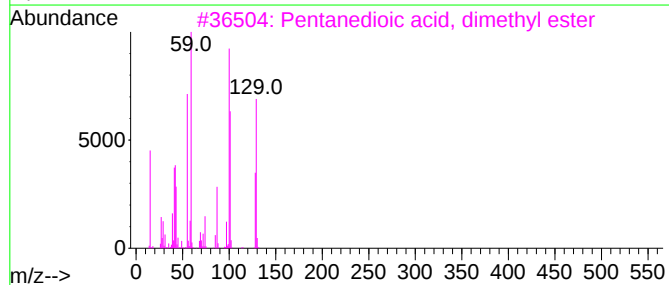
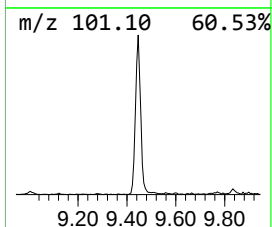
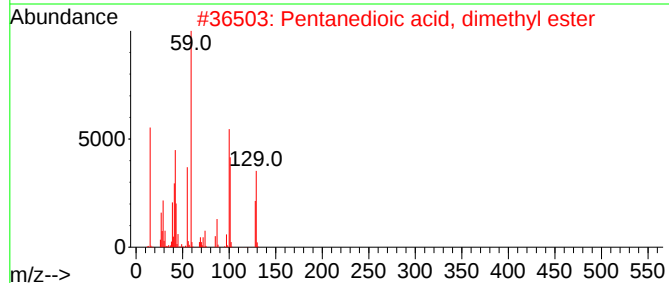
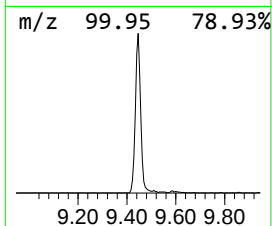
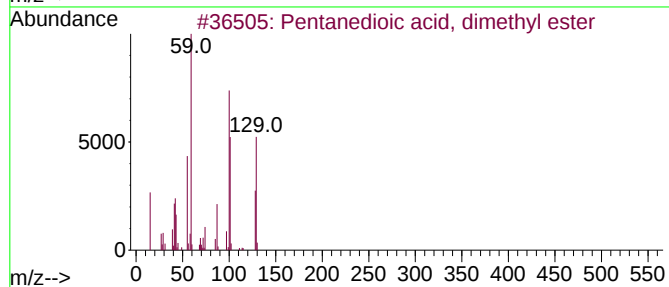
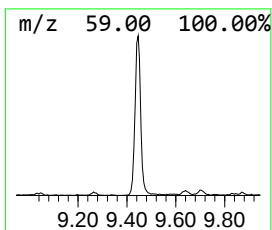
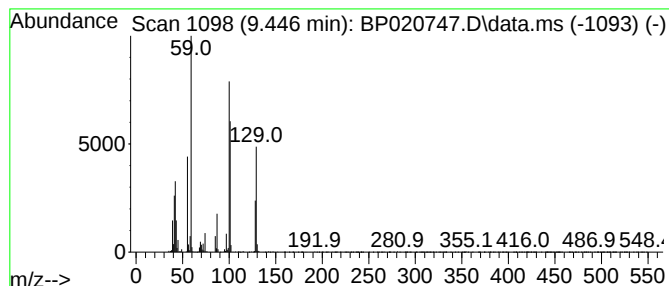
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.446	2.06 ng/ul	192343	Naphthalene-d8	10.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	90
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	72
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	72
5			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	49



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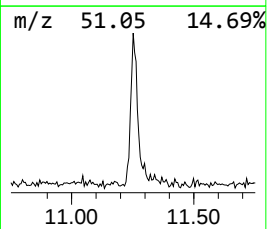
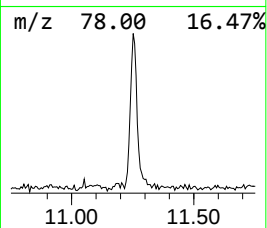
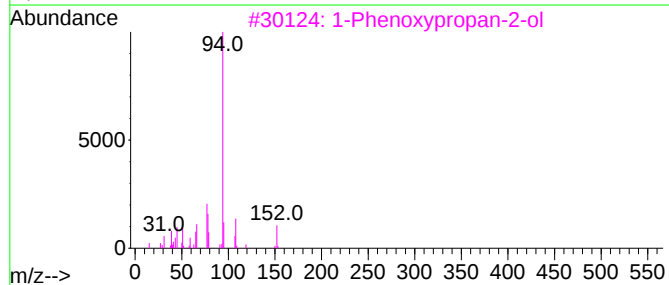
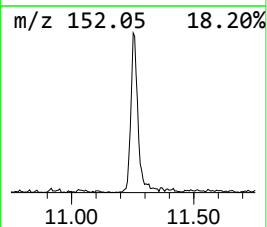
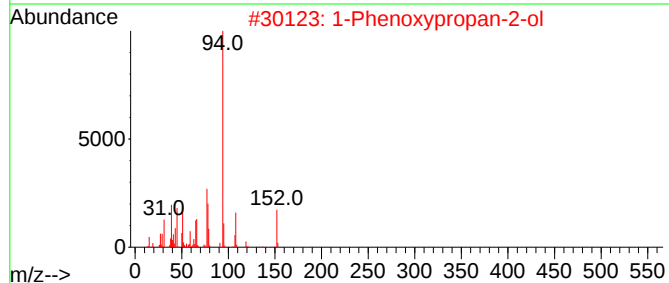
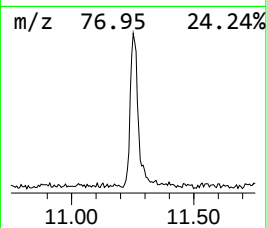
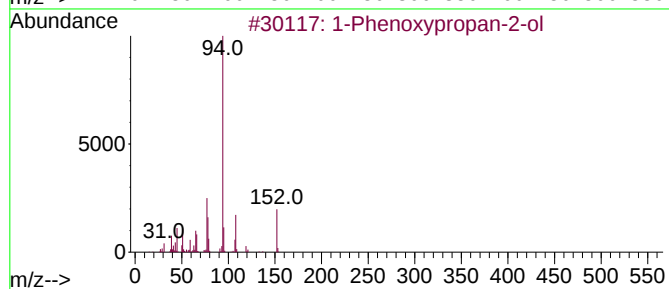
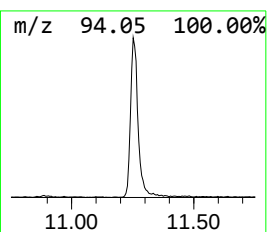
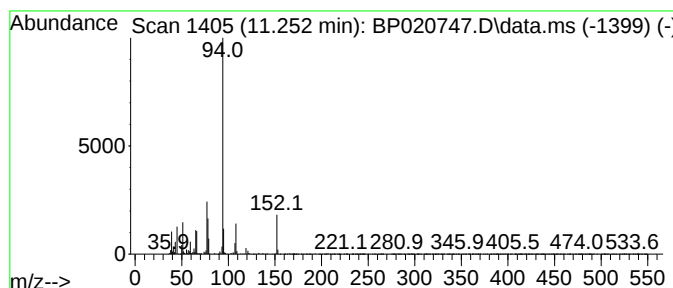
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.252	2.67 ng/ul	249292	Naphthalene-d8	10.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	96
3	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	95
4	1-Phenoxypropan-2-ol			152	C9H12O2	000770-35-4	93
5	1-Propanol, 3-phenoxy-			152	C9H12O2	006180-61-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
Operator : MA/JU
Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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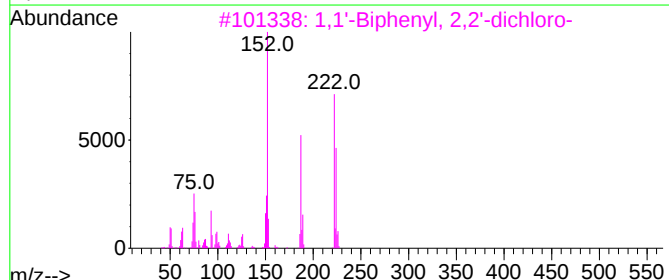
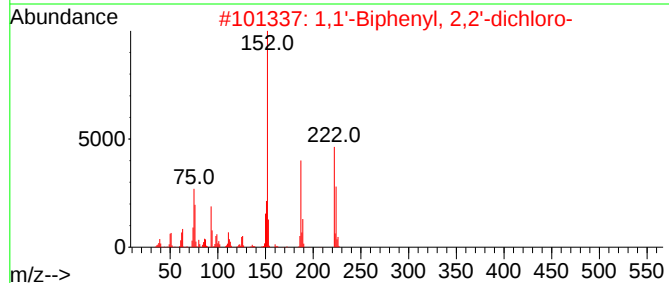
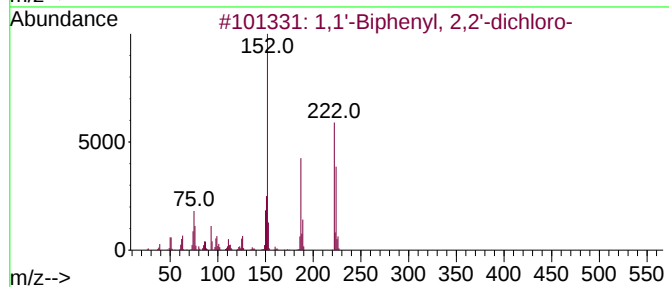
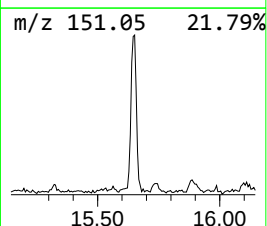
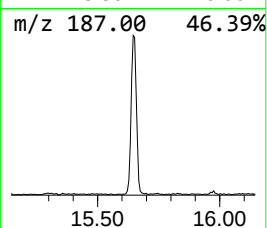
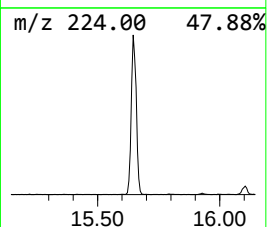
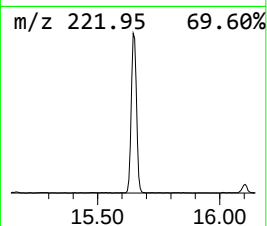
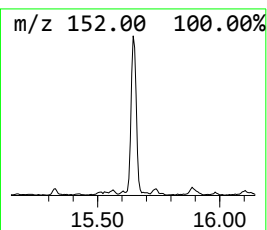
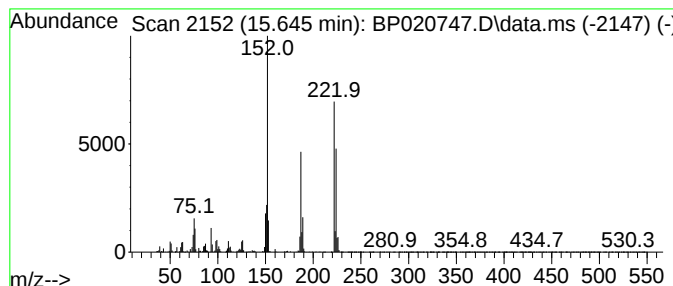
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.645	3.78 ng/ul	441575	Acenaphthene-d10	14.369

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
2			1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99
3			1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98
4			1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98
5			1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
Operator : MA/JU
Sample : P2963-02
Misc :
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Instrument :
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ClientSampleId :
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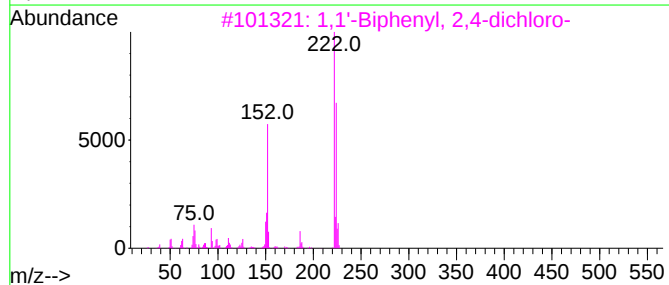
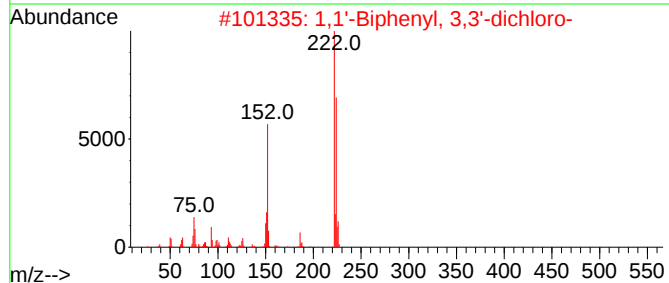
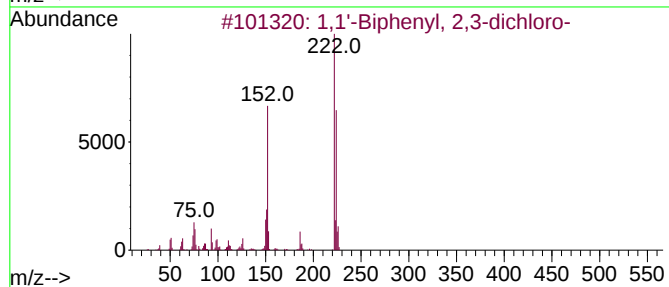
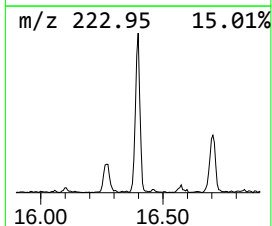
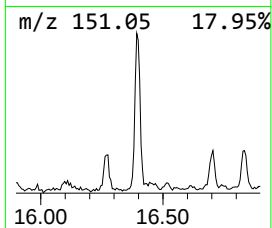
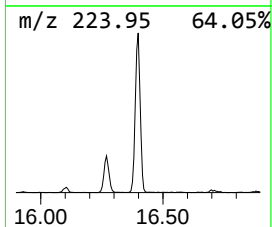
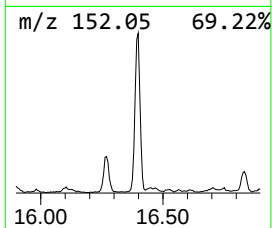
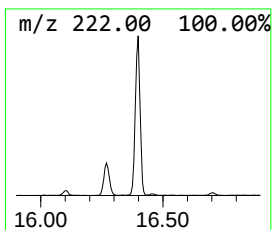
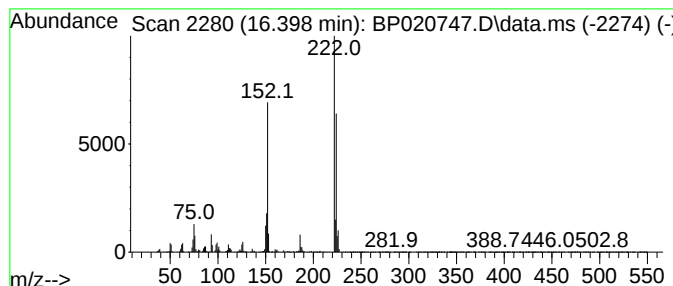
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,3-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	3.40 ng/ul	409815	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3-dichloro-	222	C12H8Cl2	016605-91-7	99		
2	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	99		
3	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	99		
4	1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	98		
5	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
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Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

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ClientSampleId :
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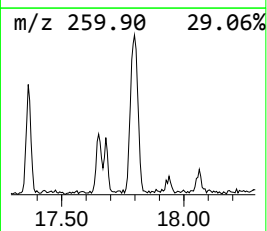
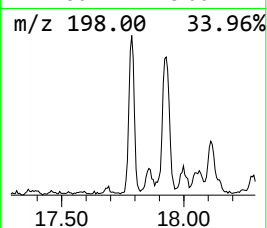
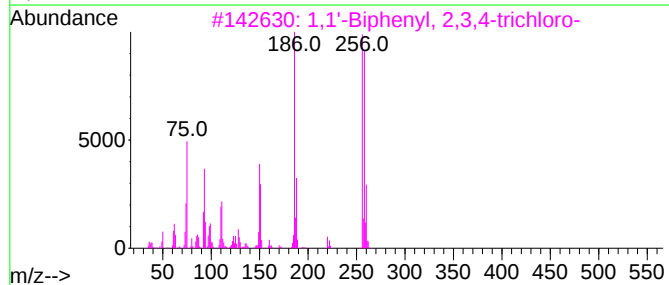
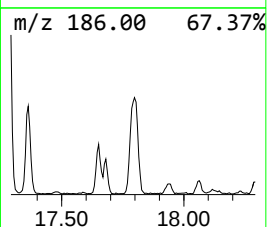
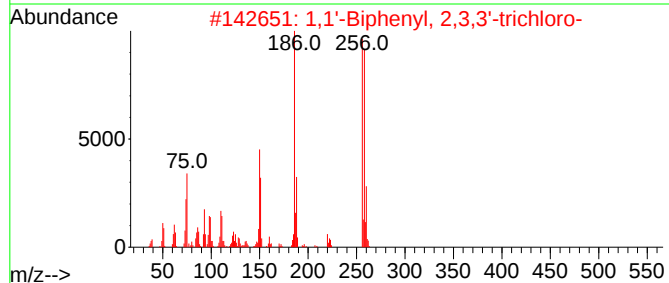
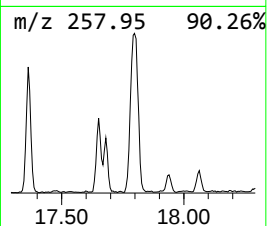
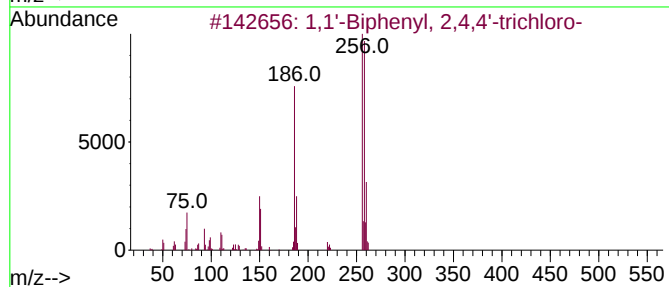
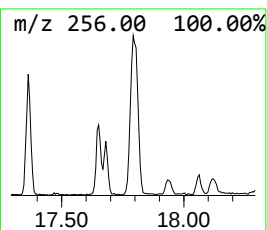
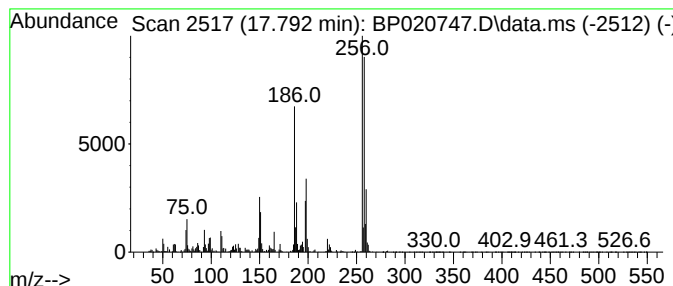
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.792	3.31 ng/ul	399008	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
3	1,1'-Biphenyl, 2,3,4-trichloro-	256	C12H7Cl3	055702-46-0	99		
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99		
5	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
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Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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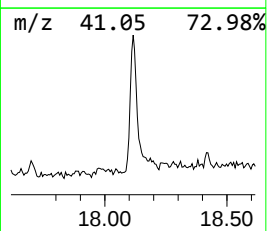
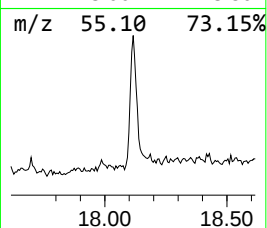
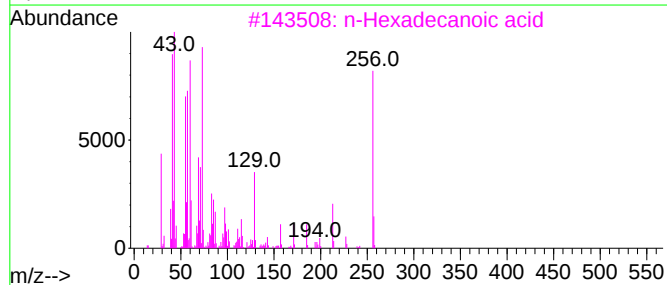
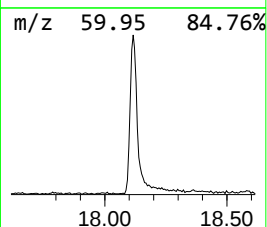
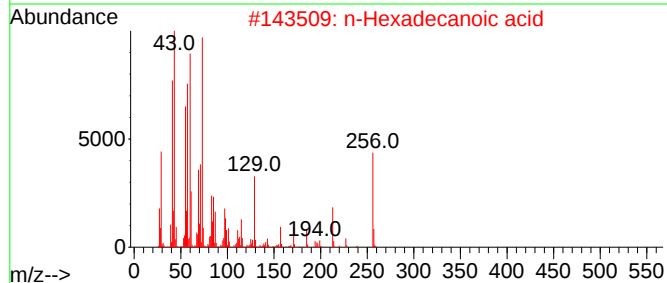
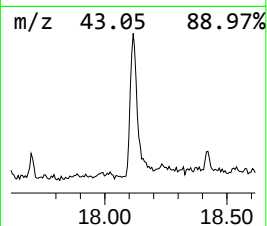
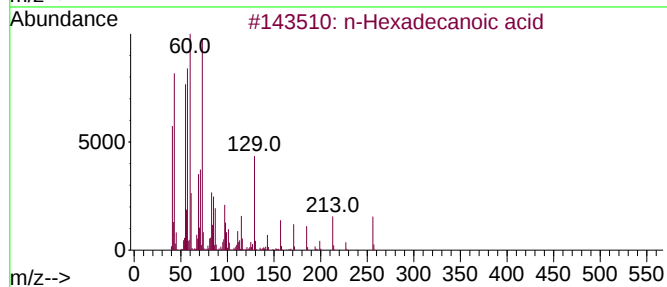
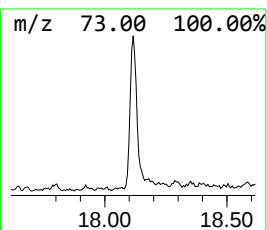
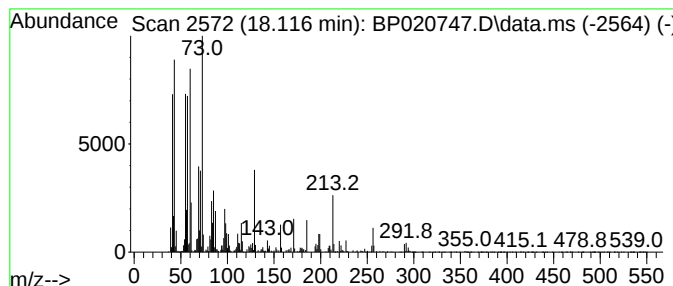
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	4.75 ng/ul	573616	Phenanthrene-d10	17.186

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	98		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
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Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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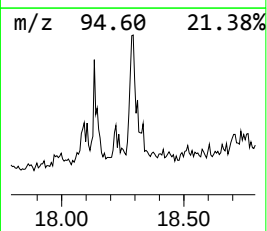
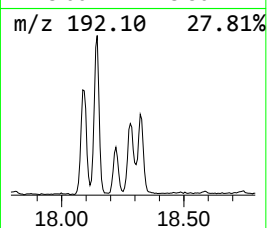
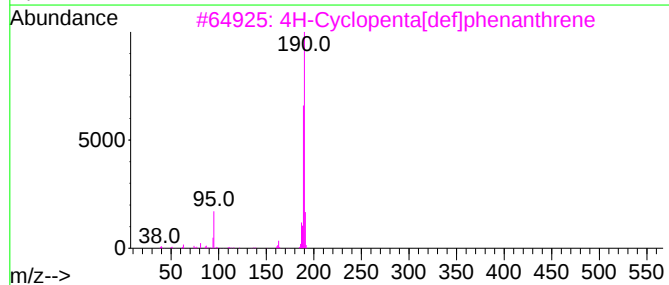
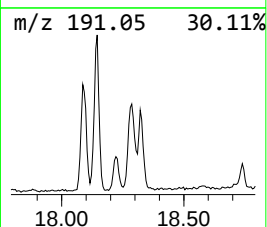
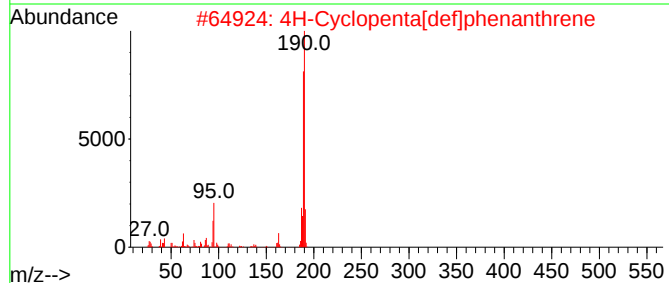
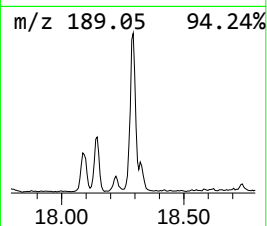
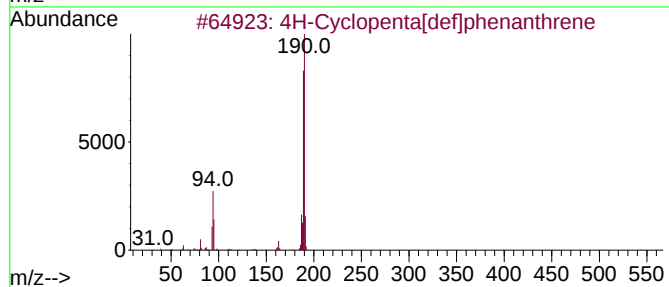
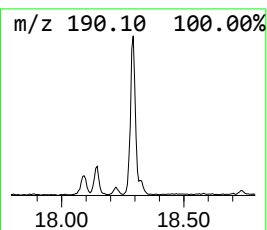
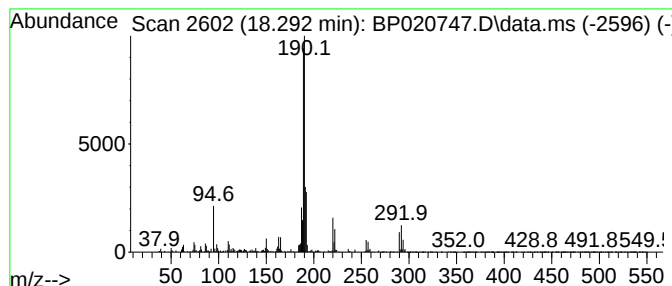
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	4.16 ng/ul	501672	Phenanthrene-d10	17.186

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4			Methyl diselenide	190	C2H6Se2	007101-31-7	46
5			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
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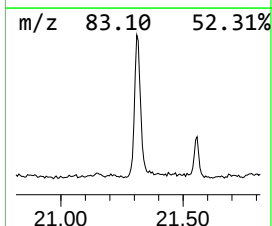
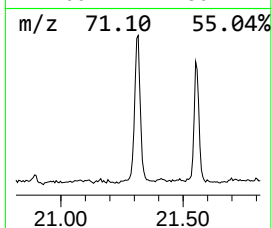
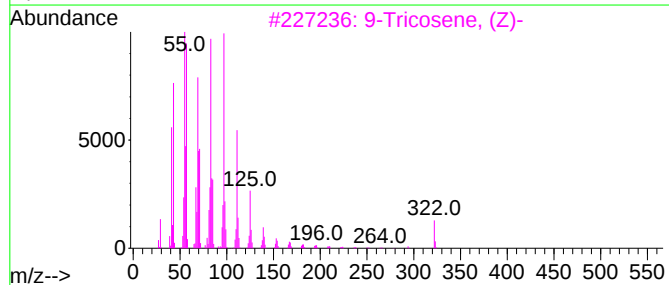
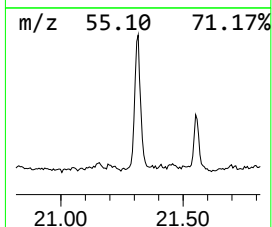
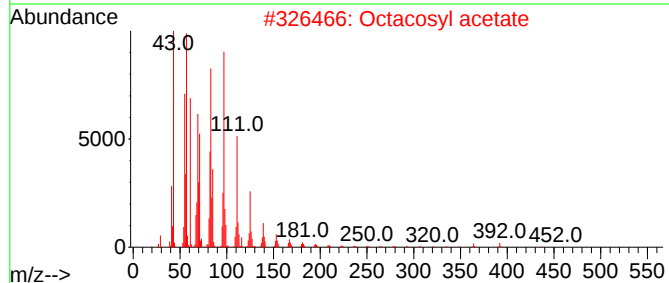
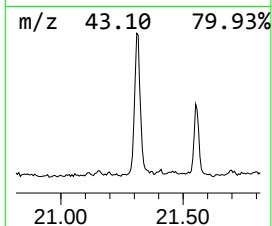
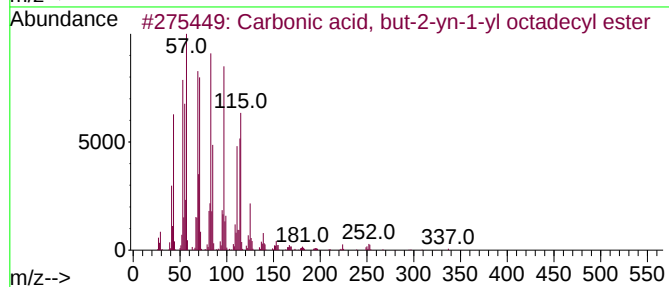
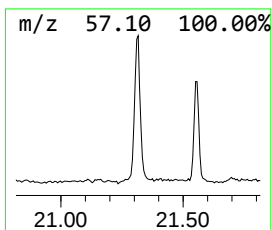
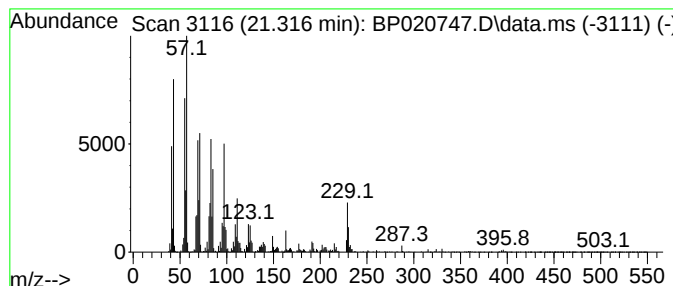
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Carbonic acid, but-2-yn-1-y... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.316	8.85 ng/ul	1515760	Chrysene-d12	21.639	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, but-2-yn-1-yl oct...	366	C23H42O3	1000383-21-3	96
2	Octacosyl acetate	452	C30H60O2	018206-97-8	95
3	9-Tricosene, (Z)-	322	C23H46	027519-02-4	90
4	1-Heneicosyl formate	340	C22H44O2	077899-03-7	86
5	Dotriacontyl heptafluorobutyrate	662	C36H65F7O2	1000351-84-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
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Misc :
ALS Vial : 37 Sample Multiplier: 1

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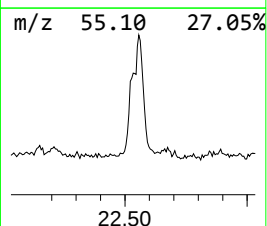
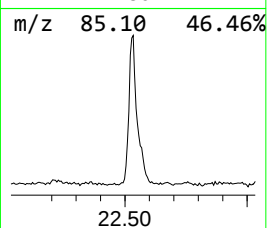
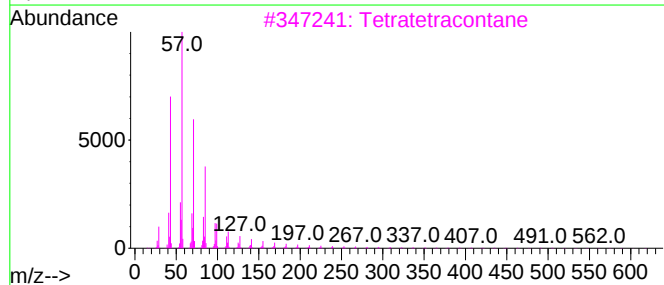
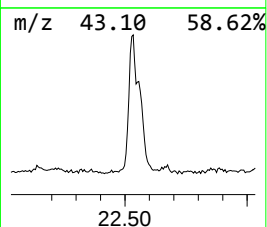
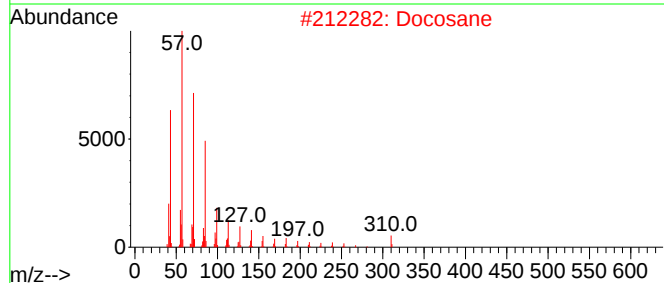
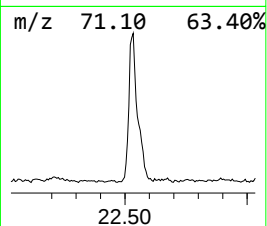
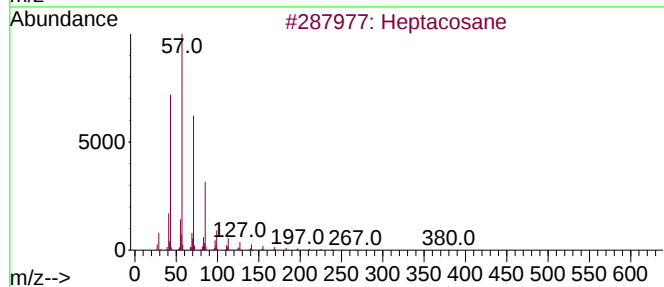
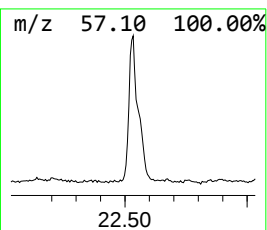
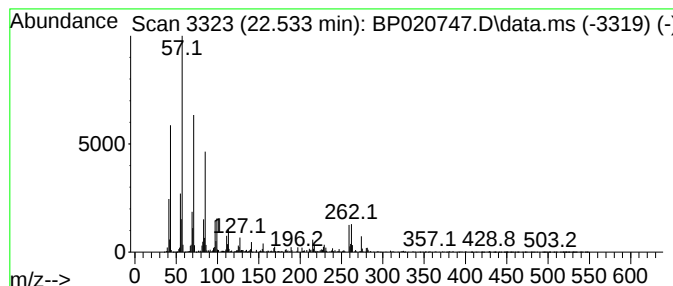
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.533	4.25 ng/ul	727611	Chrysene-d12	21.639

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	93
2			Docosane	310	C22H46	000629-97-0	89
3			Tetratetracontane	619	C44H90	007098-22-8	81
4			Hexacosane	366	C26H54	000630-01-3	81
5			2-Methylhentriacontane	451	C32H66	001720-12-3	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
Operator : MA/JU
Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4B47

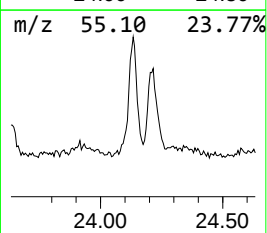
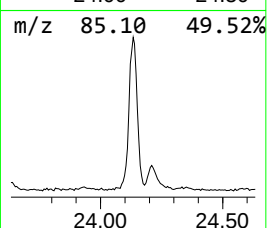
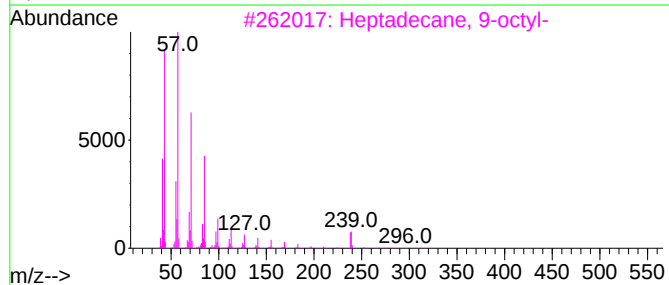
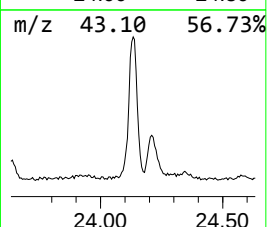
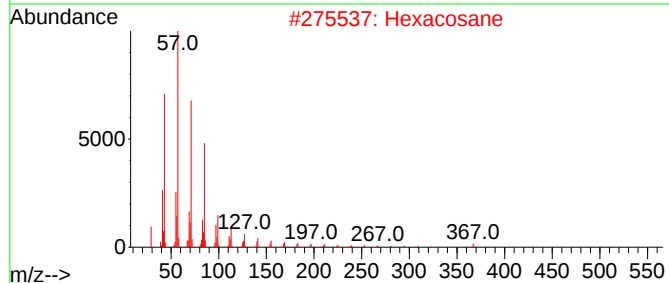
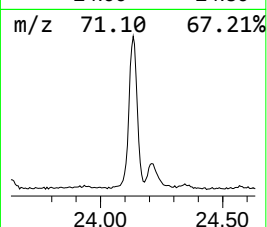
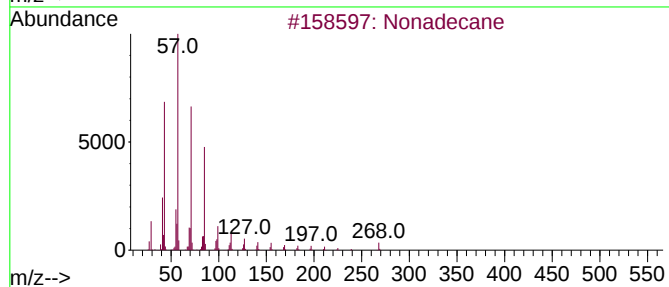
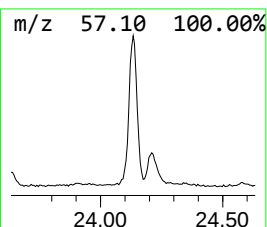
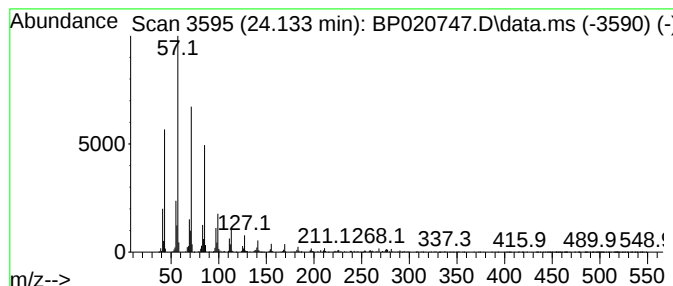
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.133	9.90 ng/ul	1241890	Perylene-d12	24.998

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	95
2		Hexacosane	366	C26H54	000630-01-3	93
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070124\
Data File : BP020747.D
Acq On : 02 Jul 2024 12:19
Operator : MA/JU
Sample : P2963-02
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4B47

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.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
unknown-01	4.893	4.3	ng/ul	287056	1	7.734	1326590	20.0
Pentanedioic ac...	9.446	2.1	ng/ul	192343	2	10.510	1867460	20.0
1-Phenoxypropan...	11.252	2.7	ng/ul	249292	2	10.510	1867460	20.0
1,1'-Biphenyl, ...	15.645	3.8	ng/ul	441575	3	14.369	2333560	20.0
1,1'-Biphenyl, ...	16.398	3.4	ng/ul	409815	4	17.186	2412740	20.0
1,1'-Biphenyl, ...	17.792	3.3	ng/ul	399008	4	17.186	2412740	20.0
n-Hexadecanoic ...	18.116	4.8	ng/ul	573616	4	17.186	2412740	20.0
4H-Cyclopenta[d...	18.292	4.2	ng/ul	501672	4	17.186	2412740	20.0
Carbonic acid, ...	21.316	8.8	ng/ul	1515760	5	21.639	3424920	20.0
(DEL) Alkane: S...	22.533	4.3	ng/ul	727611	5	21.639	3424920	20.0
(DEL) Alkane: S...	24.133	9.9	ng/ul	1241890	6	24.998	2509990	20.0