

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047883.D
 Acq On : 07 Oct 2024 13:31
 Operator : RC/JU
 Sample : P4131-21
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4E67

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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047883.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	41	45	50	rVB	32453	44763	0.20%	0.012%
2	4.528	256	262	272	rVB	37859	69793	0.31%	0.018%
3	6.181	536	543	548	rBV3	26489	61437	0.28%	0.016%
4	6.992	669	681	693	rVB2	996880	2485150	11.18%	0.648%
5	7.128	698	704	713	rBV	458725	717331	3.23%	0.187%
6	7.222	713	720	732	rVV	4455348	6603121	29.71%	1.720%
7	7.363	732	744	752	rVV2	2460354	4865772	21.89%	1.268%
8	7.798	810	818	819	rBV	780586	1027864	4.62%	0.268%
9	7.834	819	824	839	rVB	6364395	10301293	46.35%	2.684%
10	8.151	870	878	888	rBV	6335412	10334406	46.50%	2.693%
11	8.510	932	939	947	rBV	761944	1194943	5.38%	0.311%
12	8.610	947	956	968	rVB	4867405	8061677	36.27%	2.100%
13	8.881	994	1002	1008	rBV	3420604	5555618	25.00%	1.448%
14	8.951	1008	1014	1017	rVV	561446	911662	4.10%	0.238%
15	8.992	1017	1021	1030	rVB	1015968	1638235	7.37%	0.427%
16	9.363	1075	1084	1086	rBV5	21877	37599	0.17%	0.010%
17	9.528	1102	1112	1120	rBV	6037000	9866782	44.40%	2.571%
18	9.675	1130	1137	1138	rBV	465997	634994	2.86%	0.165%
19	9.704	1138	1142	1154	rVB	995806	1697005	7.64%	0.442%
20	9.828	1156	1163	1172	rVB	214325	357796	1.61%	0.093%
21	10.004	1185	1193	1203	rBV	5336654	8283411	37.27%	2.158%
22	10.239	1222	1233	1248	rVB2	3551701	7018053	31.58%	1.829%
23	10.592	1284	1293	1300	rBV	1015276	1797888	8.09%	0.468%
24	10.722	1308	1315	1322	rBV	524912	937684	4.22%	0.244%
25	10.933	1339	1351	1366	rVB	4744862	8019915	36.09%	2.090%
26	11.874	1504	1511	1519	rBV	1492759	2362212	10.63%	0.615%
27	12.016	1531	1535	1542	rVB2	34996	52871	0.24%	0.014%
28	12.145	1553	1557	1567	rVB2	58745	113749	0.51%	0.030%
29	12.251	1567	1575	1584	rBV	3140418	5073605	22.83%	1.322%
30	12.669	1641	1646	1654	rBV	80636	125302	0.56%	0.033%
31	12.863	1671	1679	1686	rBV	6435548	10206353	45.92%	2.659%
32	12.933	1686	1691	1703	rVB	1479725	2306117	10.38%	0.601%
33	13.263	1740	1747	1751	rBV	51786	70751	0.32%	0.018%
34	13.351	1757	1762	1768	rVV2	42220	62654	0.28%	0.016%
35	13.433	1771	1776	1780	rVV2	26536	40831	0.18%	0.011%
36	13.716	1818	1824	1831	rVB2	16695	38408	0.17%	0.010%

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	13.845	1839	1846	1855	rBV	1191018	1737935	7.82%	0.453%
38	14.010	1866	1874	1884	rBV	7040483	10155140	45.69%	2.646%
39	14.127	1887	1894	1896	rVV	1457567	2334685	10.50%	0.608%
40	14.151	1896	1898	1905	rVB	1227576	1661552	7.48%	0.433%
41	14.433	1939	1946	1951	rBV	1465557	2205774	9.92%	0.575%
42	14.498	1951	1957	1964	rVV	8304535	11991497	53.96%	3.124%
43	14.557	1964	1967	1973	rVB	84358	123565	0.56%	0.032%
44	14.645	1973	1982	1995	rBV2	2334634	4032062	18.14%	1.051%
45	14.751	1995	2000	2002	rVV2	123952	179256	0.81%	0.047%
46	14.798	2002	2008	2014	rVV	5002645	6948800	31.27%	1.811%
47	14.839	2014	2015	2025	rVB5	34237	54574	0.25%	0.014%
48	15.057	2043	2052	2058	rBV2	97033	155023	0.70%	0.040%
49	15.286	2087	2091	2097	rVB5	34104	47605	0.21%	0.012%
50	15.427	2108	2115	2118	rBV	1858845	2723131	12.25%	0.710%
51	15.480	2118	2124	2130	rVV2	12595861	20287103	91.28%	5.286%
52	15.539	2130	2134	2142	rVB	409602	564863	2.54%	0.147%
53	15.727	2161	2166	2171	rBV	177226	242343	1.09%	0.063%
54	15.786	2171	2176	2185	rVB	1054712	1426857	6.42%	0.372%
55	16.068	2220	2224	2229	rVB4	28459	42244	0.19%	0.011%
56	16.145	2232	2237	2244	rBV3	84475	149911	0.67%	0.039%
57	16.362	2264	2274	2282	rBV2	2145706	3269065	14.71%	0.852%
58	16.486	2288	2295	2302	rBV	10097579	15233133	68.54%	3.969%
59	16.827	2343	2353	2363	rBV	8250295	11750280	52.87%	3.062%
60	16.962	2369	2376	2386	rVB3	112941	229254	1.03%	0.060%
61	17.174	2405	2412	2415	rBV	1683703	2522072	11.35%	0.657%
62	17.215	2415	2419	2424	rVV	3599026	4954999	22.30%	1.291%
63	17.268	2424	2428	2431	rVV2	2069664	2997704	13.49%	0.781%
64	17.304	2431	2434	2445	rVB	3499931	4766122	21.45%	1.242%
65	17.580	2472	2481	2487	rBV5	38222	99058	0.45%	0.026%
66	17.780	2510	2515	2521	rVV4	99529	168791	0.76%	0.044%
67	17.845	2521	2526	2531	rVB	225099	273179	1.23%	0.071%
68	18.068	2559	2564	2570	rBV	422026	603599	2.72%	0.157%
69	18.139	2570	2576	2584	rVV	11093584	14097217	63.43%	3.673%
70	18.368	2609	2615	2621	rBV5	74028	164124	0.74%	0.043%
71	18.427	2621	2625	2631	rVB3	86021	136828	0.62%	0.036%
72	18.527	2638	2642	2646	rBV3	67368	98519	0.44%	0.026%
73	18.586	2647	2652	2659	rVB	327854	470451	2.12%	0.123%
74	18.680	2665	2668	2675	rVB9	34011	47324	0.21%	0.012%
75	18.786	2681	2686	2689	rBV5	57867	89504	0.40%	0.023%
76	18.921	2703	2709	2715	rBV2	143660	254188	1.14%	0.066%

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77	19.015	2722	2725	2728	rVB	96404	97919	0.44%	0.026%
78	19.150	2745	2748	2753	rVV2	76668	95704	0.43%	0.025%
79	19.227	2754	2761	2770	rVB	12227446	17535682	78.90%	4.569%
80	19.345	2777	2781	2789	rBV	143460	251820	1.13%	0.066%
81	19.468	2799	2802	2809	rVB	107569	144791	0.65%	0.038%
82	19.562	2812	2818	2819	rBV	2646161	3501261	15.75%	0.912%
83	19.586	2819	2822	2830	rVB	5970647	8296143	37.33%	2.162%
84	19.656	2831	2834	2839	rVV5	72451	97390	0.44%	0.025%
85	19.703	2840	2842	2847	rVB6	57679	68332	0.31%	0.018%
86	19.786	2853	2856	2863	rBV	64907	89670	0.40%	0.023%
87	20.245	2930	2934	2939	rBV2	359284	464941	2.09%	0.121%
88	20.480	2970	2974	2979	rVB	3400626	3699868	16.65%	0.964%
89	21.074	3071	3075	3086	rVB2	551039	795277	3.58%	0.207%
90	21.227	3097	3101	3107	rBV2	368582	487076	2.19%	0.127%
91	21.303	3109	3114	3121	rVV	13384432	17721295	79.74%	4.617%
92	21.386	3121	3128	3134	rVV2	10441428	18150667	81.67%	4.729%
93	21.450	3134	3139	3150	rVB	14515721	22224690	100.00%	5.791%
94	22.350	3288	3292	3302	rVB3	281927	562258	2.53%	0.146%
95	23.462	3471	3481	3490	rBV	9422448	20916180	94.11%	5.450%
96	24.191	3598	3605	3611	rBV	1183434	3078545	13.85%	0.802%
97	24.256	3611	3616	3630	rVB	1971878	4639812	20.88%	1.209%
98	24.397	3633	3640	3653	rVB	1125064	2861371	12.87%	0.746%
99	27.844	4207	4226	4246	rVB	2721770	12662450	56.97%	3.299%
100	28.838	4388	4395	4412	rVB	806670	3092527	13.91%	0.806%

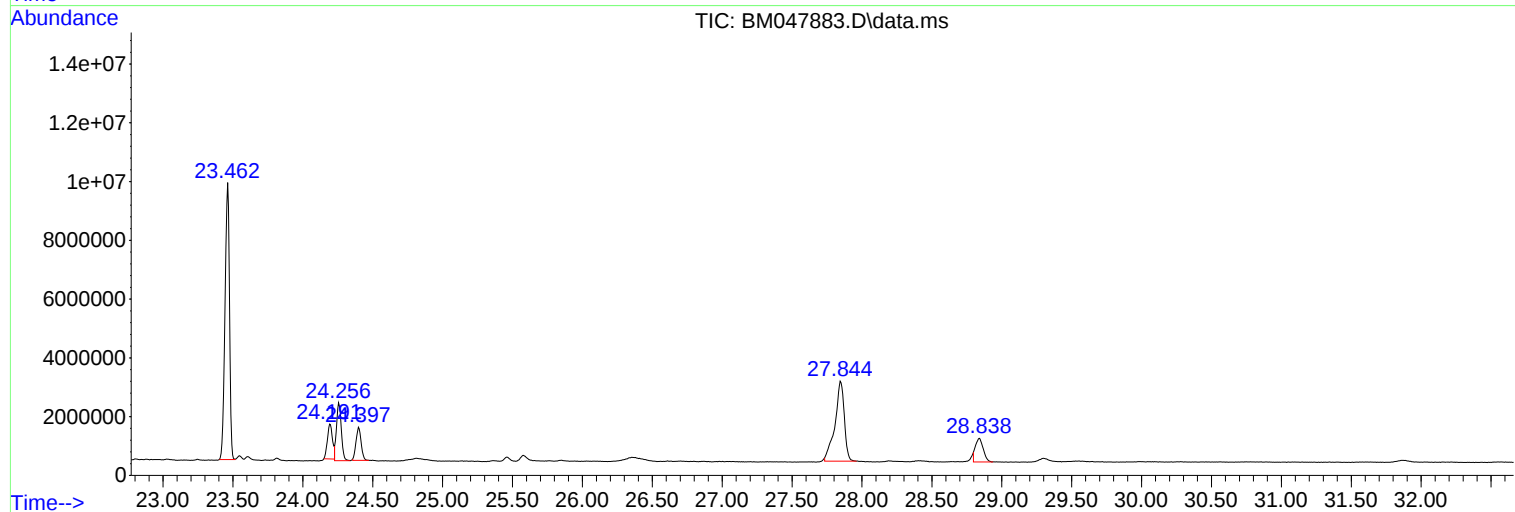
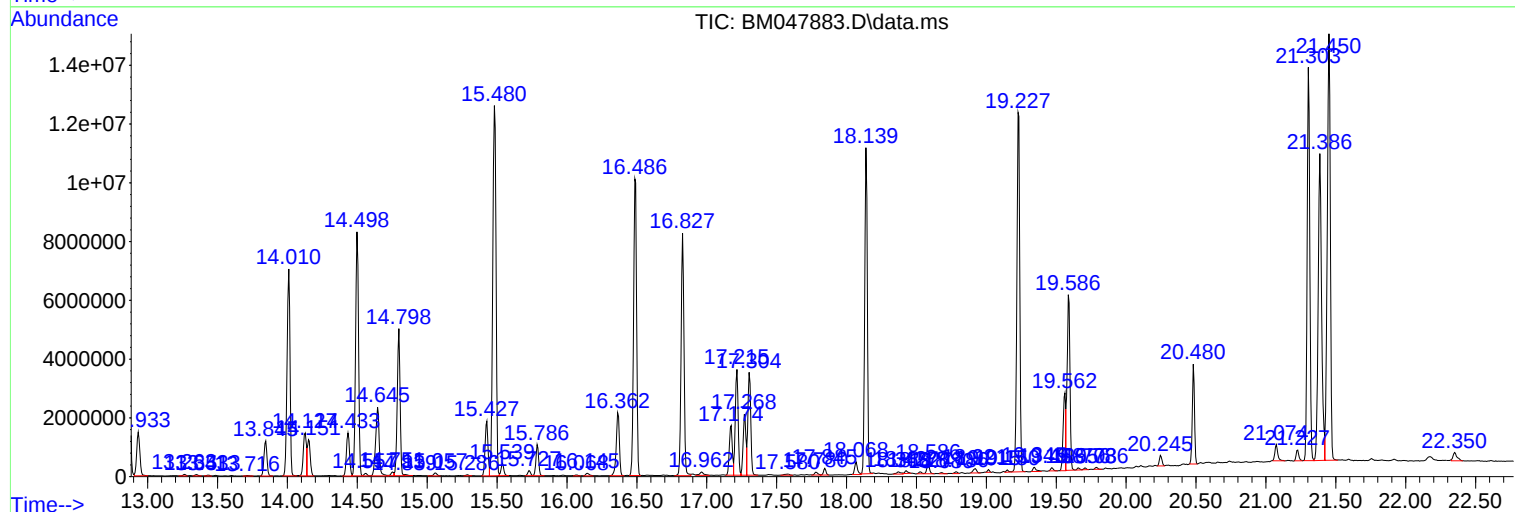
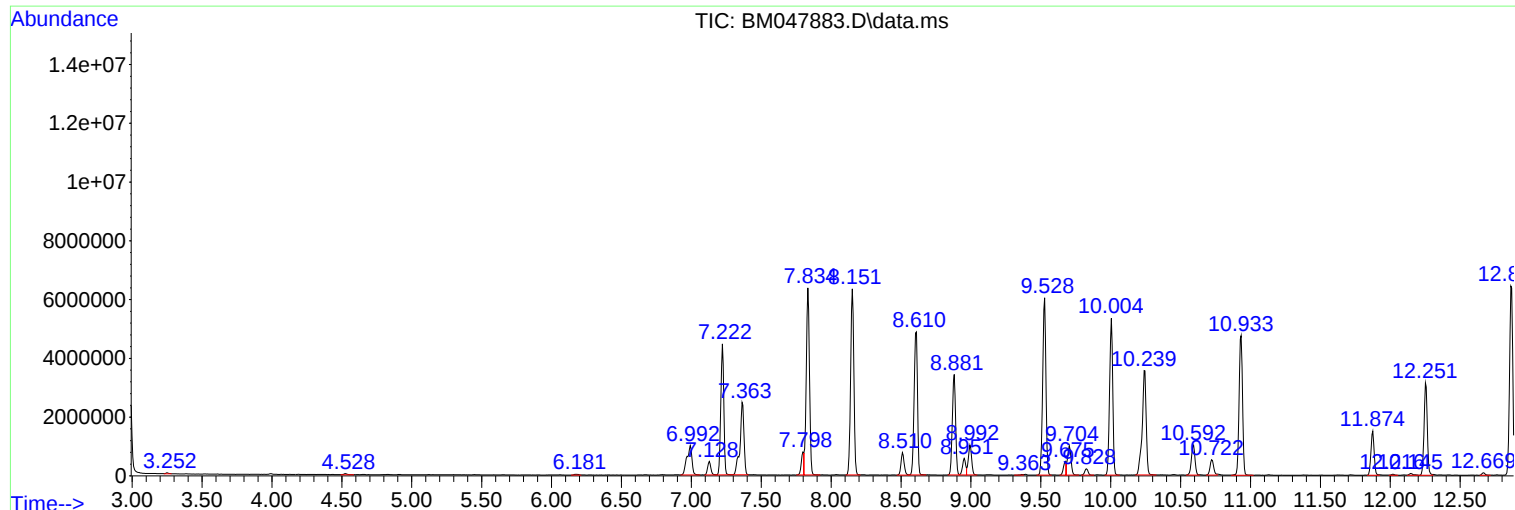
Sum of corrected areas: 383802045

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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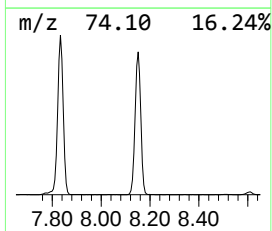
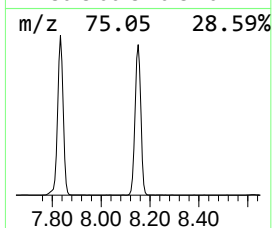
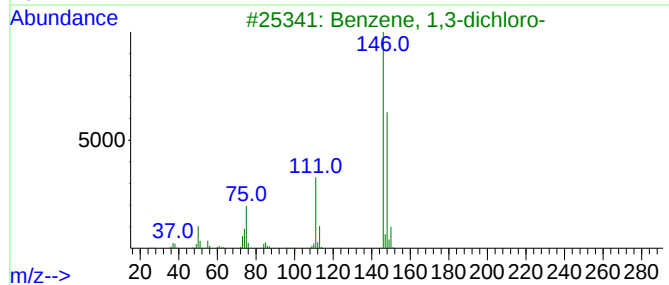
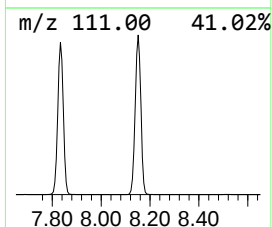
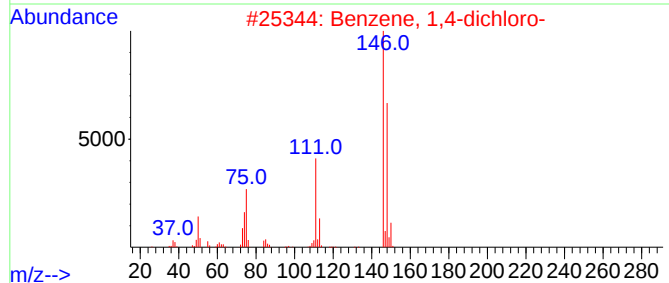
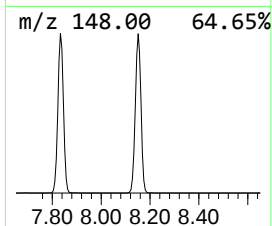
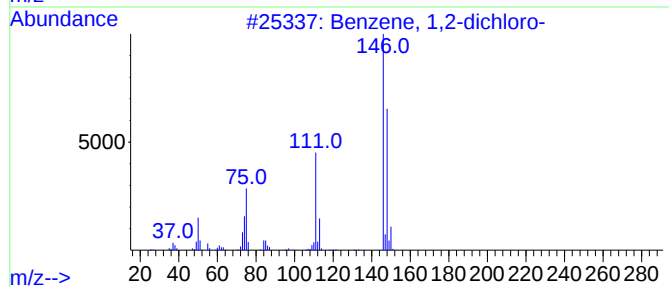
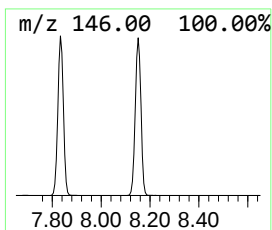
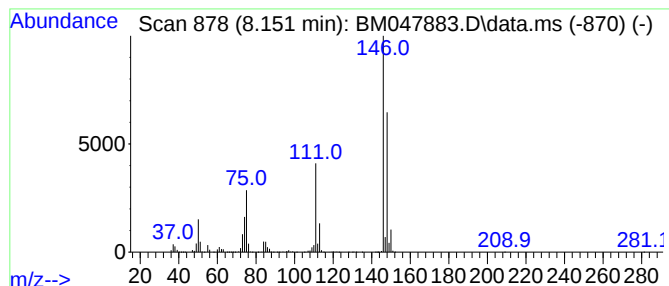
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,2-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.151	201.09 ng/ul	10334400	1,4-Dichlorobenzene-d4	7.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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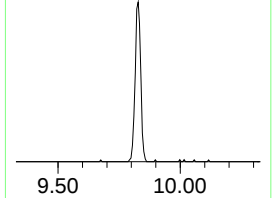
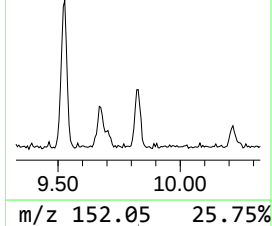
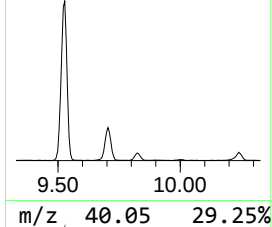
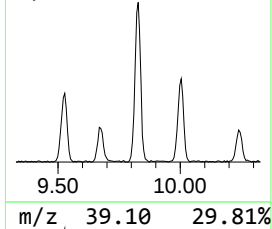
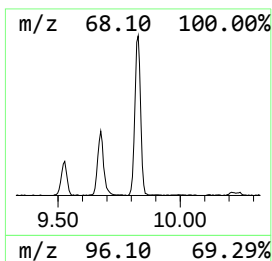
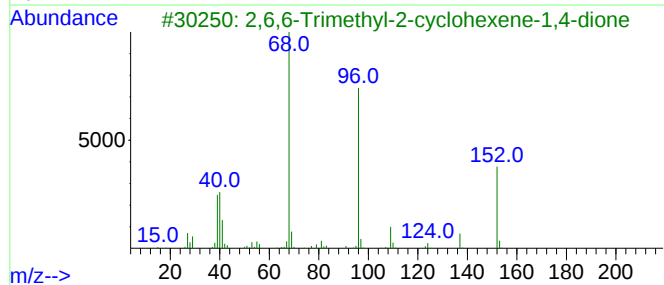
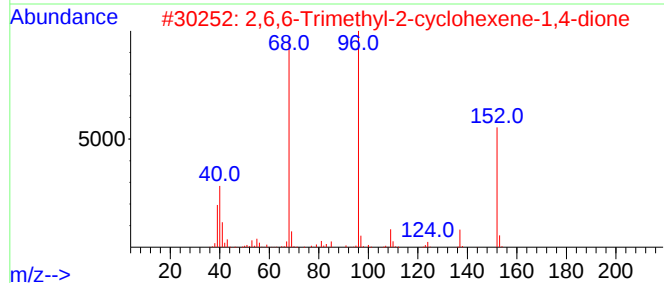
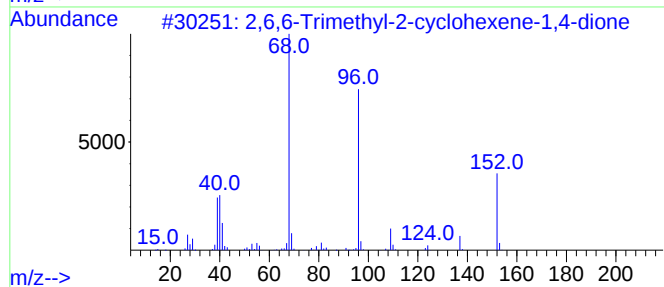
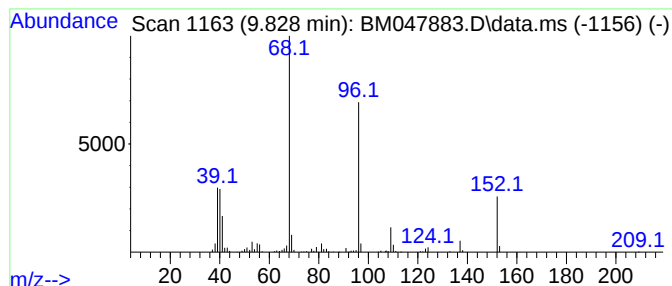
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.828	3.98 ng/ul	357796	Naphthalene-d8	10.592

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	95		
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
4	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	76		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53		



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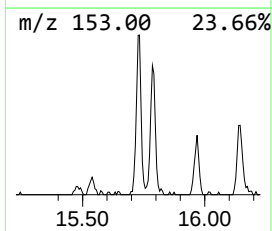
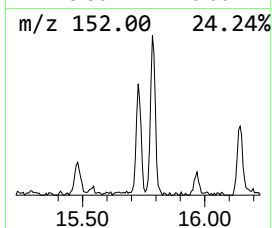
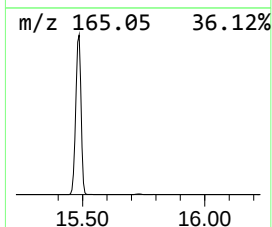
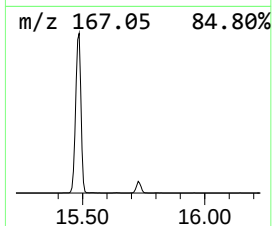
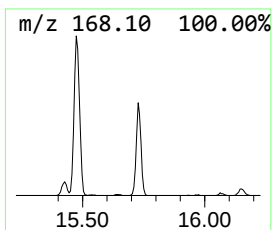
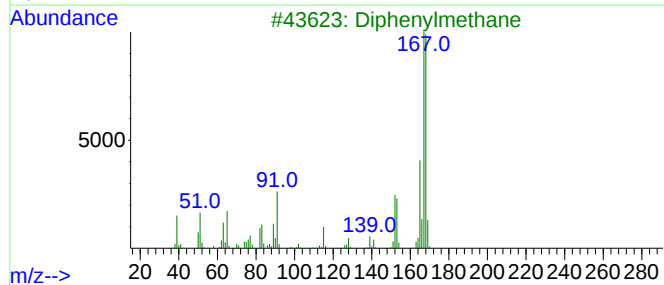
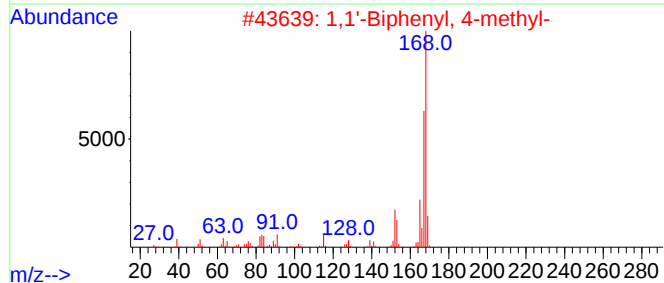
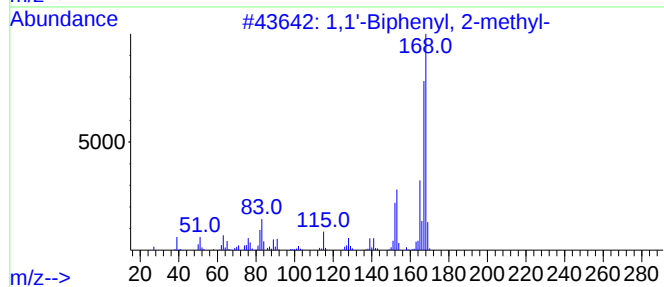
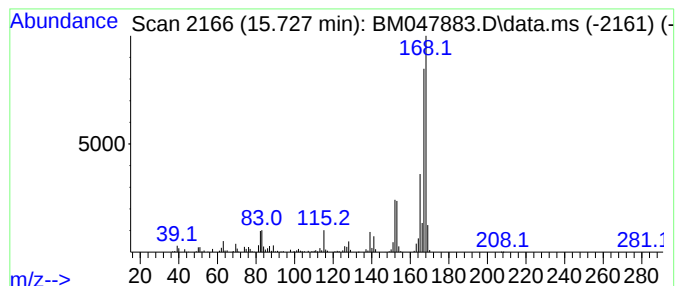
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.727	2.20 ng/ul	242343	Acenaphthene-d10	14.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-			168	C13H12	000643-58-3	95
2	1,1'-Biphenyl, 4-methyl-			168	C13H12	000644-08-6	95
3	Diphenylmethane			168	C13H12	000101-81-5	93
4	Diphenylmethane			168	C13H12	000101-81-5	93
5	1,1'-Biphenyl, 4-methyl-			168	C13H12	000644-08-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047883.D
Acq On : 07 Oct 2024 13:31
Operator : RC/JU
Sample : P4131-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4E67

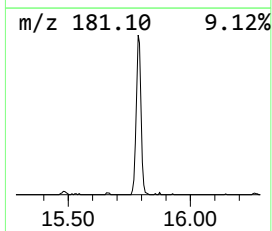
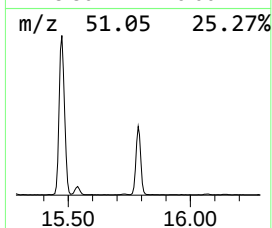
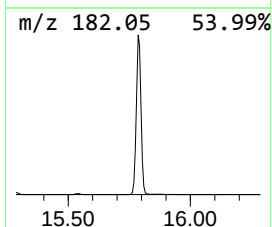
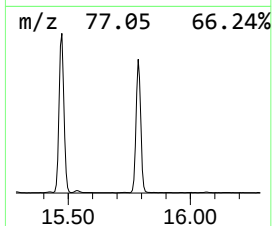
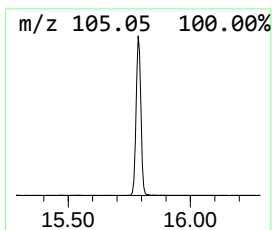
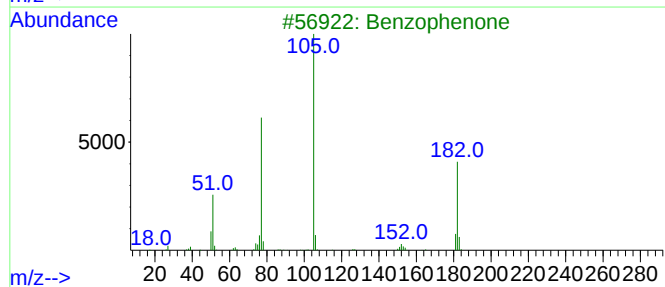
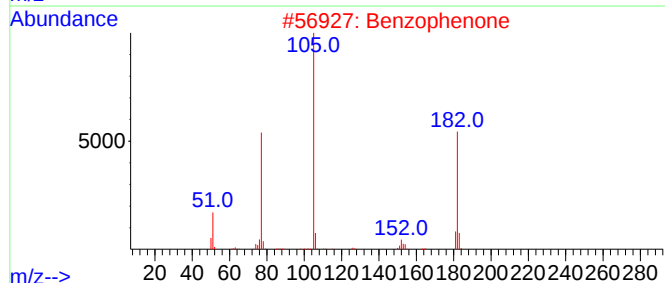
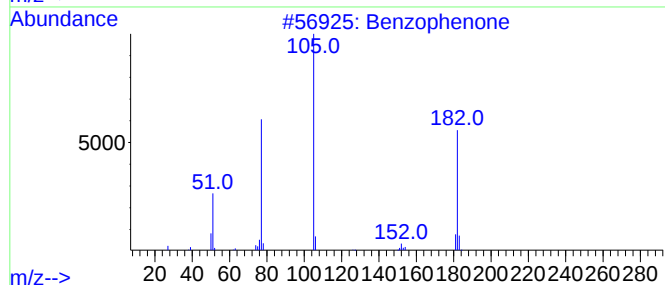
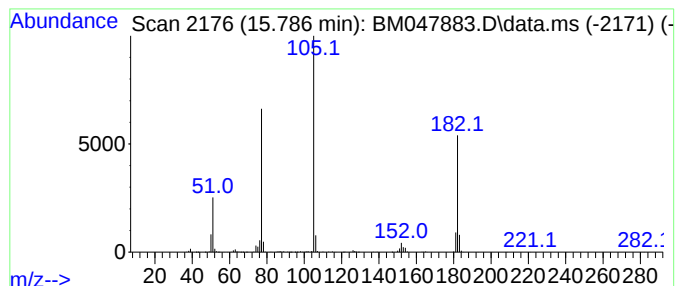
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.786	12.94 ng/ul	1426860	Acenaphthene-d10	14.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Azobenzene	182	C12H10N2	000103-33-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047883.D
Acq On : 07 Oct 2024 13:31
Operator : RC/JU
Sample : P4131-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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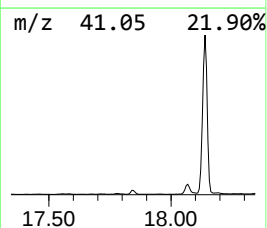
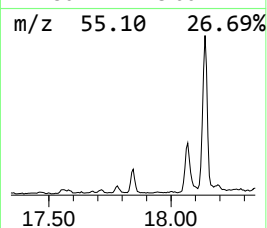
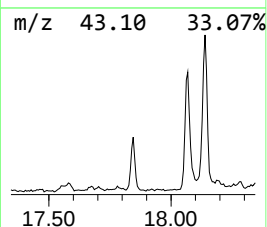
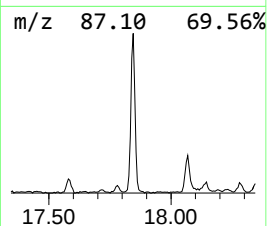
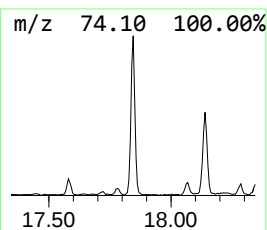
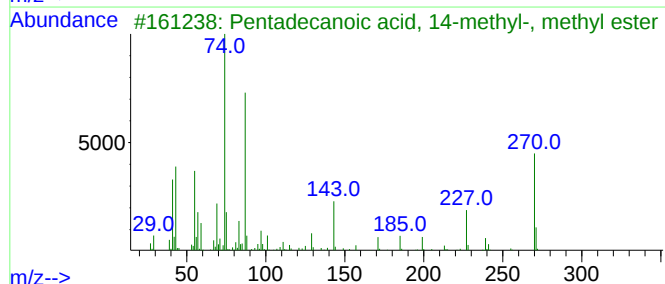
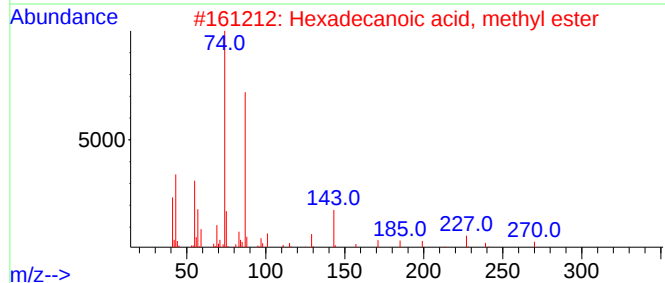
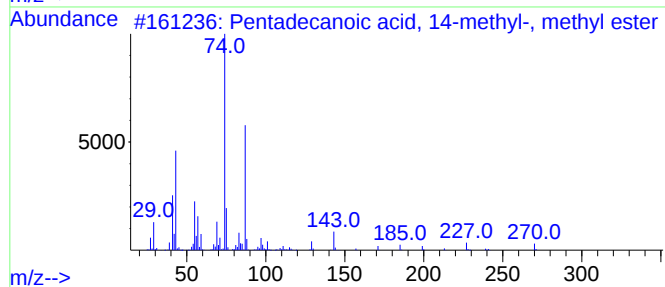
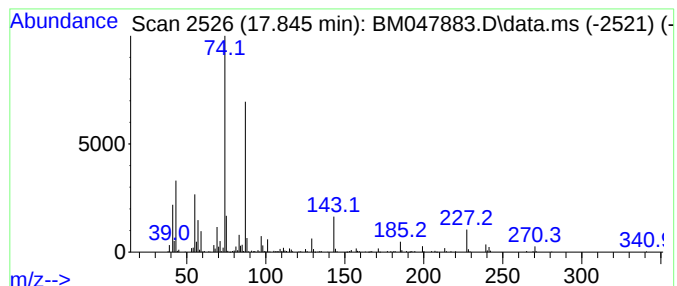
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecanoic acid, 14-meth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.845	2.17 ng/ul	273179	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
2			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
5			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047883.D
Acq On : 07 Oct 2024 13:31
Operator : RC/JU
Sample : P4131-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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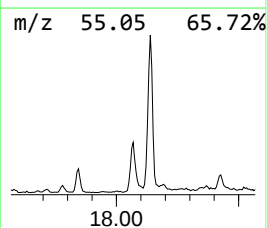
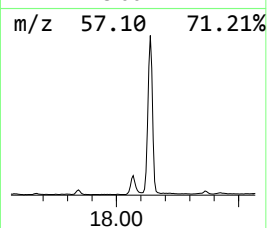
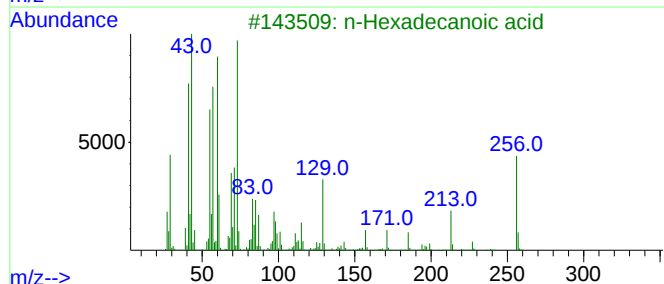
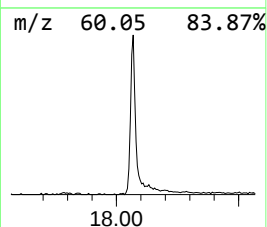
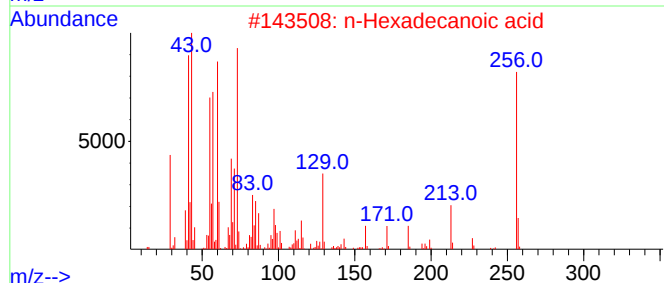
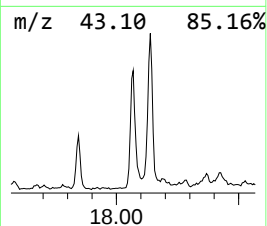
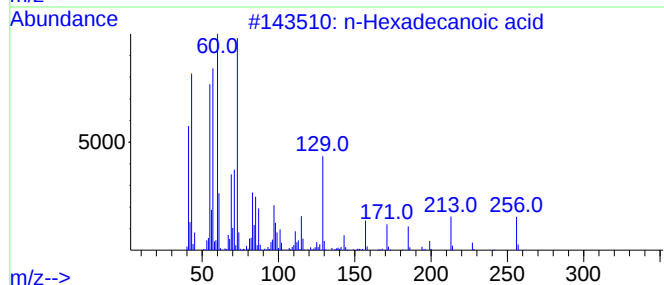
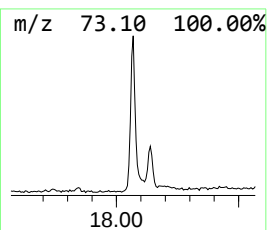
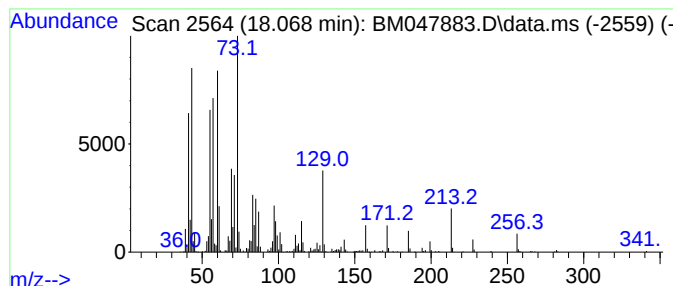
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.068	4.79 ng/ul	603599	Phenanthrene-d10	17.174

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047883.D
Acq On : 07 Oct 2024 13:31
Operator : RC/JU
Sample : P4131-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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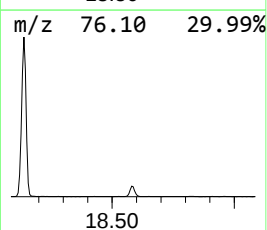
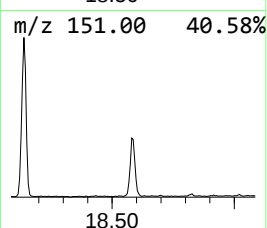
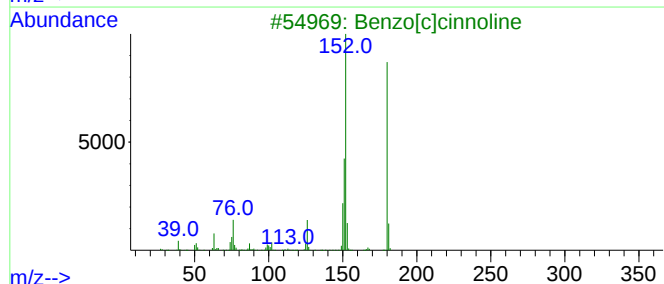
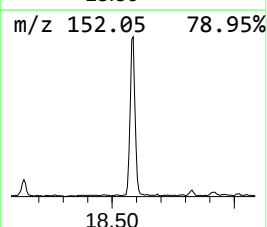
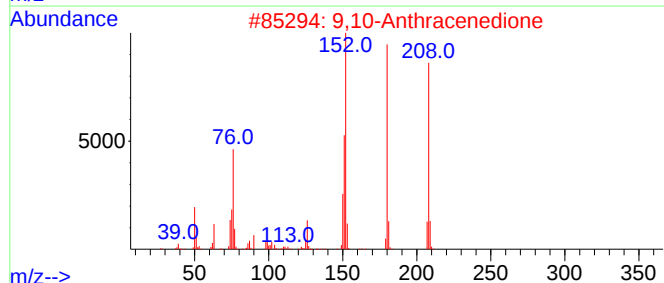
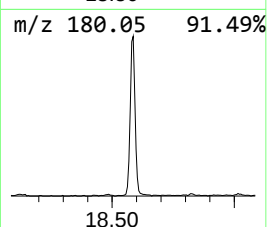
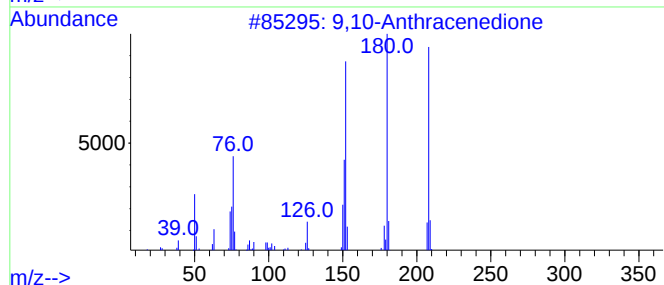
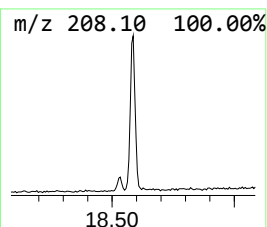
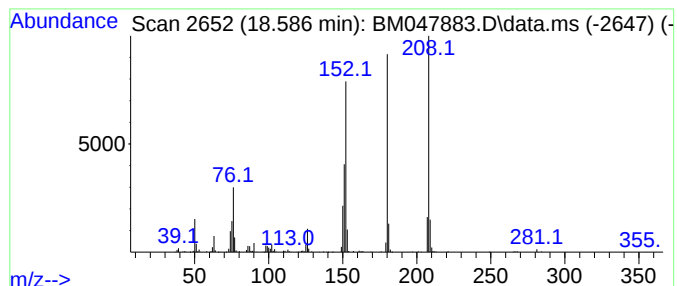
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.586	3.73 ng/ul	470451	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	95
5	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
Data File : BM047883.D
Acq On : 07 Oct 2024 13:31
Operator : RC/JU
Sample : P4131-21
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
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2,6,6-Trimethyl...	9.828	4.0	ng/ul	357796	2	10.592	1797890	20.0
1,1'-Biphenyl, ...	15.727	2.2	ng/ul	242343	3	14.433	2205770	20.0
Benzophenone	15.786	12.9	ng/ul	1426860	3	14.433	2205770	20.0
Pentadecanoic a...	17.845	2.2	ng/ul	273179	4	17.174	2522070	20.0
n-Hexadecanoic ...	18.068	4.8	ng/ul	603599	4	17.174	2522070	20.0
9,10-Anthracene...	18.586	3.7	ng/ul	470451	4	17.174	2522070	20.0
10,18-Bisnorabi...	18.921	2.0	ng/ul	254188	4	17.174	2522070	20.0