

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
 Data File : BM048075.D
 Acq On : 16 Oct 2024 00:22
 Operator : RC/JU
 Sample : P4350-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EOAJ9

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: BM048075.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.046	6	10	24	rVB	1563074	2630664	39.65%	3.654%
2	3.240	40	43	50	rVB	37543	53525	0.81%	0.074%
3	3.669	112	116	121	rBV2	36583	59843	0.90%	0.083%
4	3.722	122	125	127	rVV3	17080	23382	0.35%	0.032%
5	3.757	127	131	137	rVB	52229	70410	1.06%	0.098%
6	3.975	163	168	176	rVV5	24418	52053	0.78%	0.072%
7	4.351	227	232	236	rBV3	23567	37547	0.57%	0.052%
8	4.445	242	248	251	rBV	22598	39915	0.60%	0.055%
9	4.510	254	259	261	rBV	71607	113896	1.72%	0.158%
10	4.575	267	270	275	rVB	48955	67547	1.02%	0.094%
11	4.775	300	304	309	rBV6	11090	16981	0.26%	0.024%
12	4.893	319	324	330	rVB	37682	61433	0.93%	0.085%
13	5.587	438	442	447	rBV6	15585	23526	0.35%	0.033%
14	5.787	470	476	482	rBV3	21497	42384	0.64%	0.059%
15	6.040	514	519	526	rBV5	15107	28134	0.42%	0.039%
16	6.163	530	540	549	rVB2	65314	161794	2.44%	0.225%
17	6.416	568	583	590	rVV	83470	174016	2.62%	0.242%
18	6.598	607	614	616	rBV3	15507	27537	0.42%	0.038%
19	6.639	616	621	628	rVB4	27103	47686	0.72%	0.066%
20	6.957	665	675	690	rBV2	515421	1017979	15.34%	1.414%
21	7.110	695	701	709	rBV	442035	684550	10.32%	0.951%
22	7.316	729	736	747	rBV	571724	938637	14.15%	1.304%
23	7.692	795	800	807	rBV8	12628	29885	0.45%	0.042%
24	7.781	807	815	821	rBV	1031491	1646507	24.82%	2.287%
25	7.851	823	827	834	rVB3	21564	41190	0.62%	0.057%
26	8.033	851	858	863	rBV3	8363	20071	0.30%	0.028%
27	8.092	863	868	874	rVV	48029	73968	1.11%	0.103%
28	8.492	926	936	948	rBV	583983	1020142	15.38%	1.417%
29	8.598	949	954	963	rVV	183771	315456	4.75%	0.438%
30	8.933	1004	1011	1021	rBV	475064	801222	12.08%	1.113%
31	9.363	1078	1084	1091	rVB2	34692	59809	0.90%	0.083%
32	9.498	1100	1107	1115	rBV2	49036	83724	1.26%	0.116%
33	9.657	1124	1134	1148	rBV	389714	697890	10.52%	0.969%
34	9.833	1157	1164	1171	rBV	29257	71869	1.08%	0.100%
35	9.910	1171	1177	1184	rVB8	8933	22761	0.34%	0.032%
36	10.016	1189	1195	1199	rBV3	16925	29165	0.44%	0.041%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
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37	10.198	1219	1226	1236	rBV	579478	1108312	16.71%	1.539%
38	10.275	1237	1239	1245	rVB4	22917	40193	0.61%	0.056%
39	10.569	1280	1289	1297	rBV	1367096	2346559	35.37%	3.259%
40	10.710	1305	1313	1319	rBV	222138	468655	7.06%	0.651%
41	11.027	1362	1367	1374	rVB10	11227	21628	0.33%	0.030%
42	11.122	1376	1383	1392	rBV10	20226	49529	0.75%	0.069%
43	11.780	1487	1495	1501	rBV10	10041	27297	0.41%	0.038%
44	11.857	1501	1508	1512	rVV6	9962	18973	0.29%	0.026%
45	12.163	1551	1560	1566	rVV5	12198	31008	0.47%	0.043%
46	12.233	1569	1572	1581	rVB2	12944	23653	0.36%	0.033%
47	12.786	1658	1666	1670	rBV7	10048	22314	0.34%	0.031%
48	13.245	1740	1744	1751	rVB6	17449	30336	0.46%	0.042%
49	13.586	1795	1802	1805	rBV9	8004	19979	0.30%	0.028%
50	13.739	1824	1828	1833	rVB7	14175	24683	0.37%	0.034%
51	13.827	1833	1843	1853	rBV	998033	1516062	22.85%	2.106%
52	14.110	1880	1891	1902	rBV	1265462	1931512	29.11%	2.683%
53	14.415	1935	1943	1950	rBV	1977080	2992399	45.10%	4.156%
54	14.480	1950	1954	1961	rVB	113527	172889	2.61%	0.240%
55	14.551	1962	1966	1974	rVB6	15038	27752	0.42%	0.039%
56	14.633	1974	1980	1992	rBV	219765	478524	7.21%	0.665%
57	14.780	2001	2005	2007	rBV3	18494	29157	0.44%	0.040%
58	14.821	2007	2012	2021	rVB2	59061	145771	2.20%	0.202%
59	15.086	2054	2057	2060	rVB5	14318	18782	0.28%	0.026%
60	15.210	2073	2078	2081	rBV4	21281	36950	0.56%	0.051%
61	15.409	2106	2112	2118	rVV	1644233	2417678	36.44%	3.358%
62	15.462	2118	2121	2127	rVV2	138846	201692	3.04%	0.280%
63	15.533	2128	2133	2140	rVV	221533	327112	4.93%	0.454%
64	15.627	2146	2149	2154	rBV6	17899	28826	0.43%	0.040%
65	15.774	2168	2174	2181	rVB2	270497	405148	6.11%	0.563%
66	15.904	2191	2196	2204	rBV2	28590	70303	1.06%	0.098%
67	16.139	2232	2236	2243	rBV7	27199	59238	0.89%	0.082%
68	16.368	2271	2275	2280	rBV3	33983	53175	0.80%	0.074%
69	16.515	2297	2300	2304	rVB5	22212	33281	0.50%	0.046%
70	16.562	2305	2308	2314	rBV7	30134	44645	0.67%	0.062%
71	16.815	2347	2351	2355	rBV	88869	122548	1.85%	0.170%
72	16.974	2374	2378	2386	rVB	118930	199065	3.00%	0.276%
73	17.156	2403	2409	2413	rBV	2302153	3401560	51.27%	4.724%
74	17.198	2413	2416	2421	rVV	2200730	3094928	46.65%	4.299%
75	17.256	2421	2426	2430	rVV	1793170	2626915	39.59%	3.648%
76	17.292	2430	2432	2437	rVV	363436	407441	6.14%	0.566%

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77	17.351	2439	2442	2447	rVB	45628	70023	1.06%	0.097%
78	17.562	2474	2478	2483	rVB	140343	194939	2.94%	0.271%
79	17.927	2536	2540	2547	rBV6	59075	135324	2.04%	0.188%
80	18.050	2554	2561	2565	rBV2	745673	1342137	20.23%	1.864%
81	18.227	2586	2591	2595	rBV2	386609	627753	9.46%	0.872%
82	18.350	2609	2612	2615	rBV2	42262	51891	0.78%	0.072%
83	18.533	2639	2643	2646	rBV	176835	231668	3.49%	0.322%
84	18.574	2646	2650	2655	rVV2	306539	437136	6.59%	0.607%
85	18.621	2656	2658	2662	rVB	61254	71221	1.07%	0.099%
86	19.092	2734	2738	2746	rBV2	186560	389359	5.87%	0.541%
87	19.215	2753	2759	2766	rVB	4824370	6634604	100.00%	9.215%
88	19.327	2774	2778	2787	rBV3	343954	532460	8.03%	0.740%
89	19.544	2810	2815	2818	rBV3	2278972	3634598	54.78%	5.048%
90	19.574	2818	2820	2828	rVB	2684398	3421913	51.58%	4.753%
91	19.756	2848	2851	2854	rVB	129748	152577	2.30%	0.212%
92	19.897	2871	2875	2877	rBV2	149124	249147	3.76%	0.346%
93	20.068	2901	2904	2910	rVB2	470860	627369	9.46%	0.871%
94	20.168	2918	2921	2925	rVB	296678	355039	5.35%	0.493%
95	20.539	2982	2984	2988	rVB	266077	240230	3.62%	0.334%
96	20.815	3028	3031	3038	rVB	264178	381311	5.75%	0.530%
97	21.386	3121	3128	3132	rBV2	2699711	6283272	94.70%	8.727%
98	21.433	3132	3136	3147	rVB	1678693	2768518	41.73%	3.845%
99	23.438	3471	3477	3484	rBV	1170671	3186806	48.03%	4.426%
100	24.385	3631	3638	3646	rVB	1429959	3617039	54.52%	5.024%

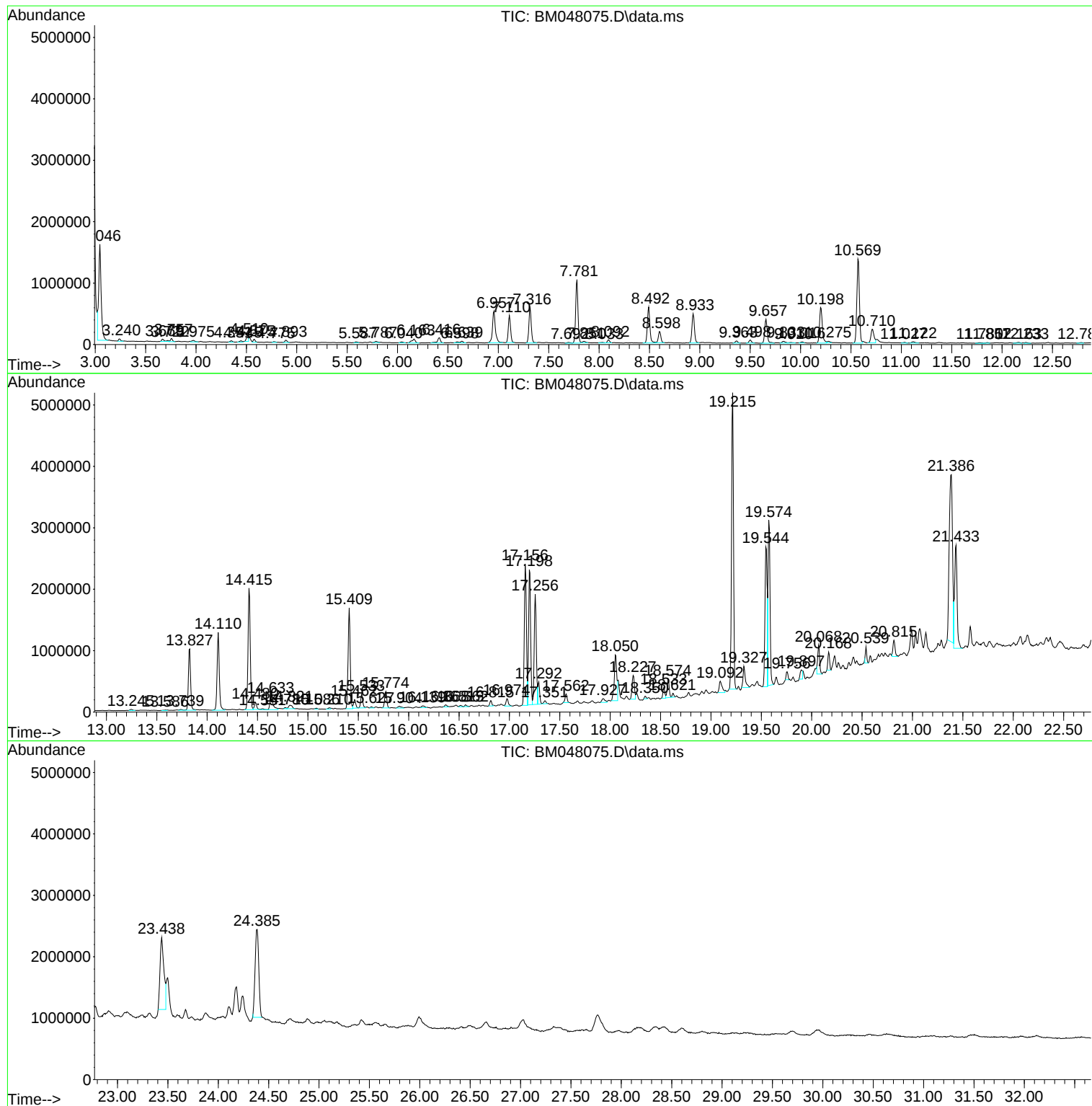
Sum of corrected areas: 71999904

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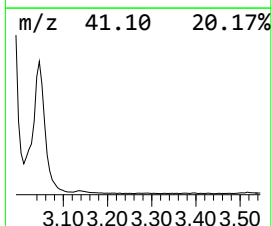
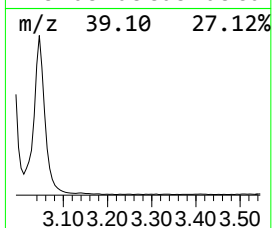
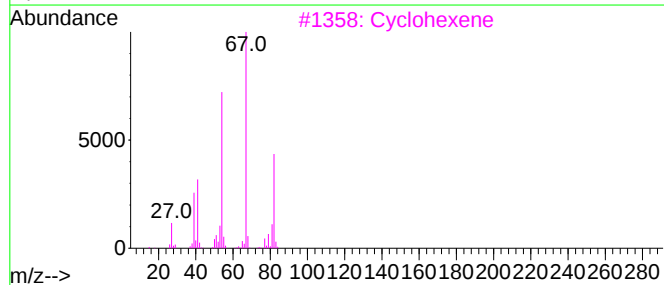
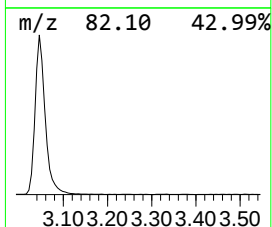
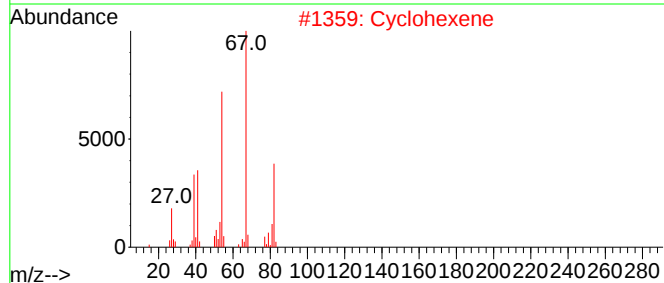
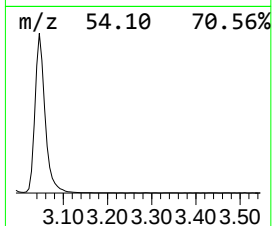
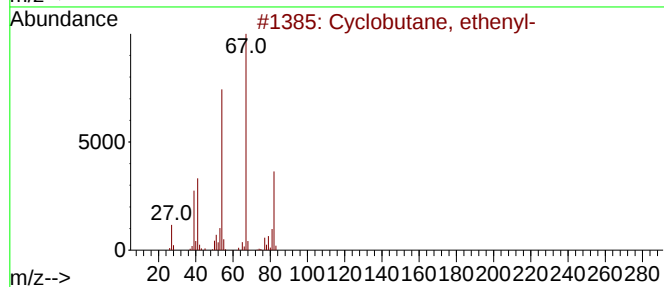
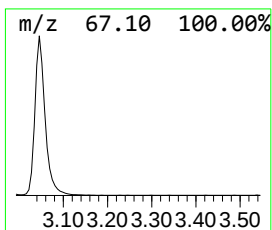
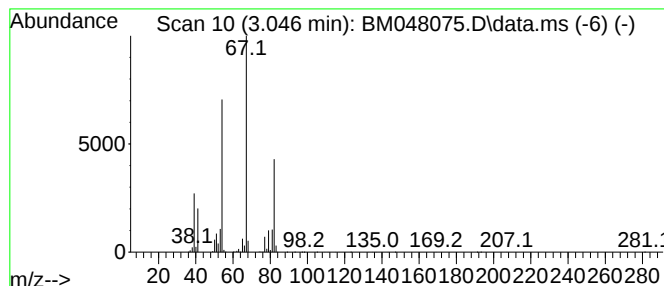
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclobutane, ethenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.046	31.95 ng/ul	2630660	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutane, ethenyl-	82	C6H10	002597-49-1	93
2			Cyclohexene	82	C6H10	000110-83-8	90
3			Cyclohexene	82	C6H10	000110-83-8	90
4			Cyclohexene	82	C6H10	000110-83-8	90
5			Ethylidenecyclobutane	82	C6H10	001528-21-8	87



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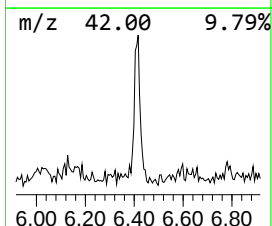
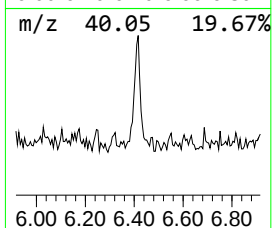
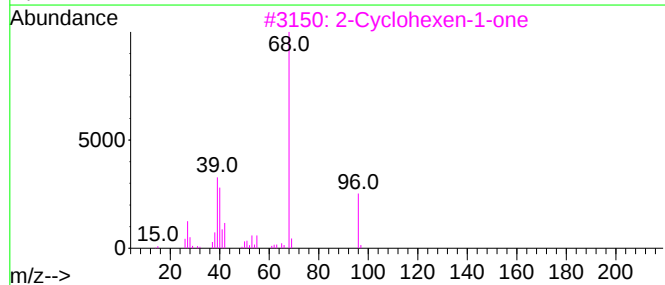
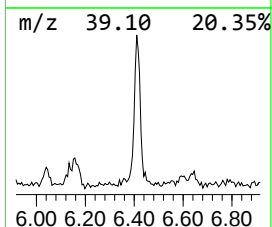
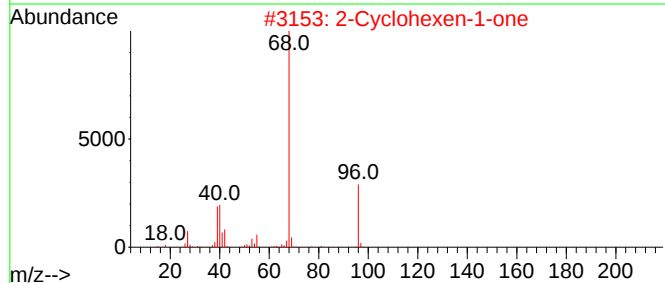
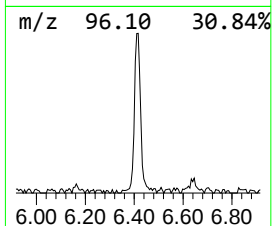
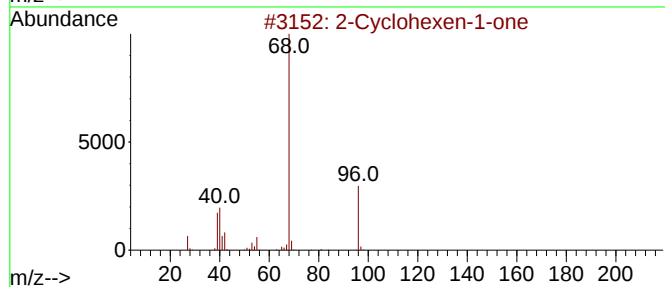
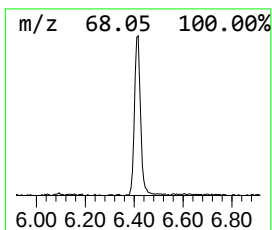
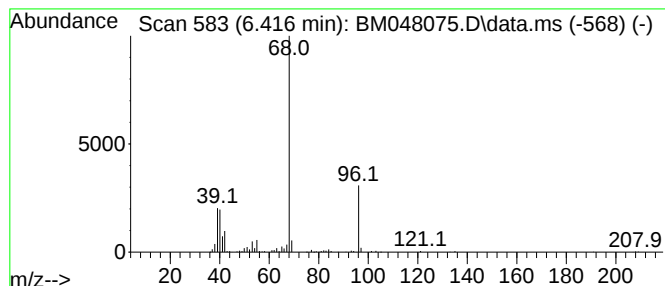
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Cyclohexen-1-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.416	2.11 ng/ul	174016	1,4-Dichlorobenzene-d4	7.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	91
2	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	90
3	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	86
4	2-Cyclohexen-1-one			96	C6H8O	000930-68-7	49
5	3-Butyn-2-amine, 2-methyl-			83	C5H9N	002978-58-7	45



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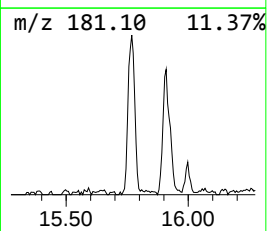
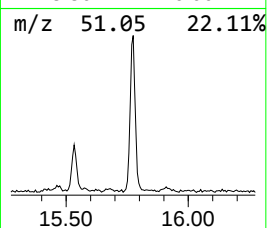
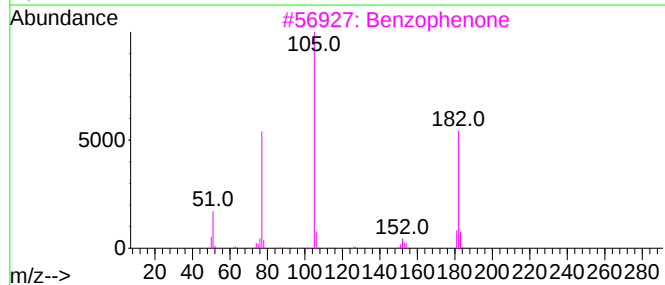
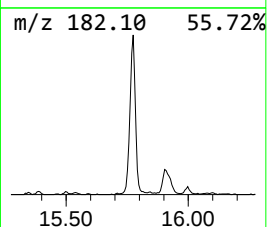
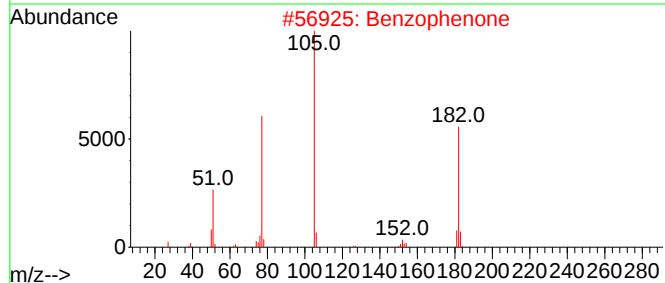
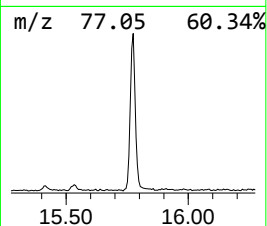
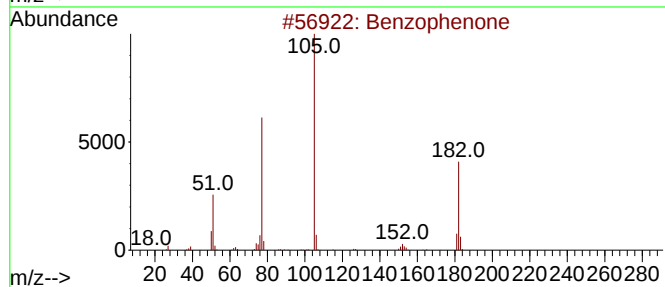
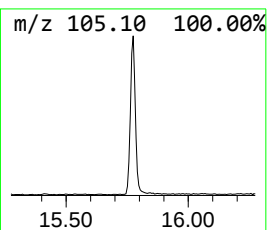
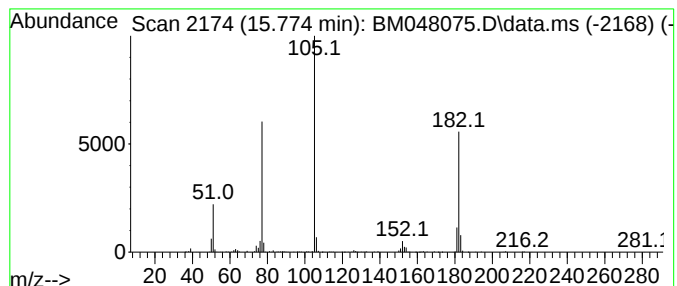
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.774	2.71 ng/ul	405148	Acenaphthene-d10	14.415

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048075.D
Acq On : 16 Oct 2024 00:22
Operator : RC/JU
Sample : P4350-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAJ9

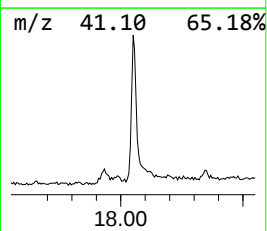
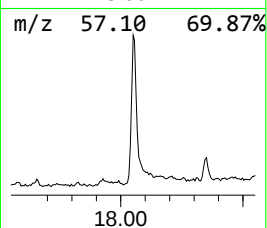
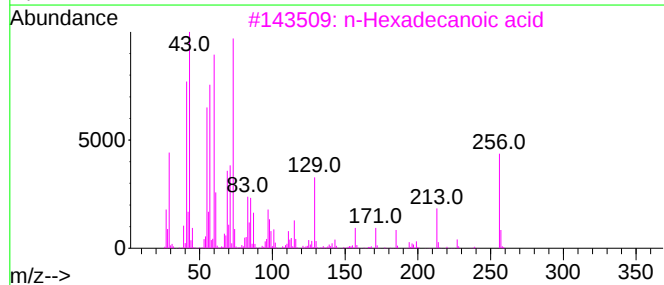
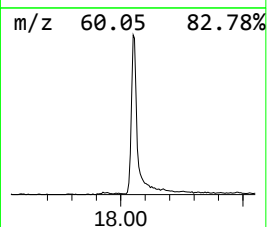
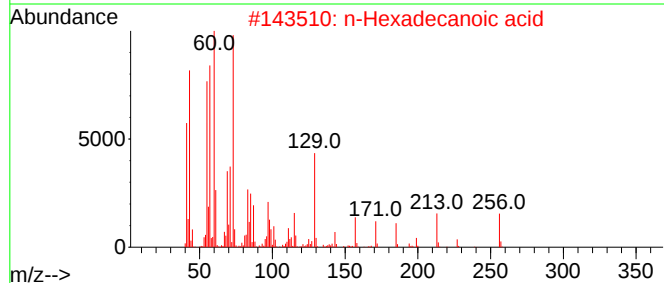
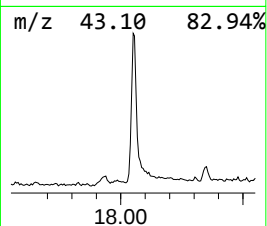
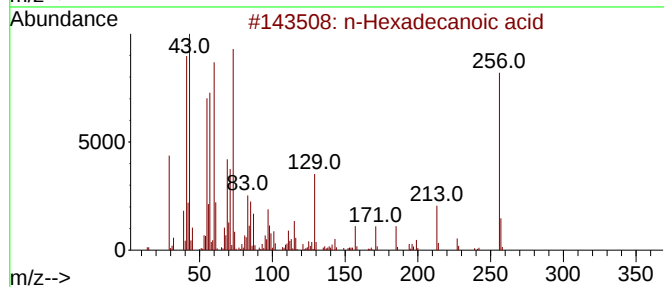
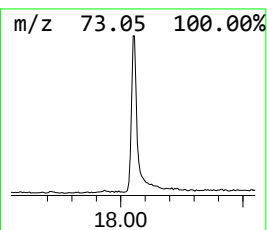
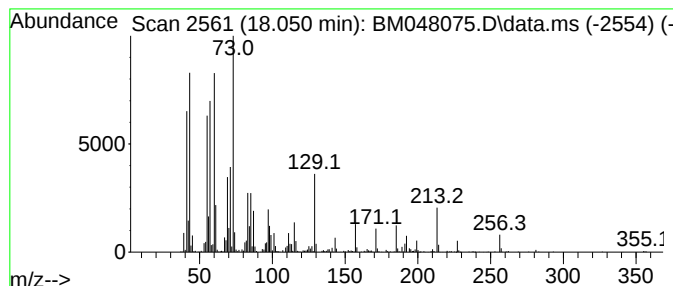
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.051	7.89 ng/ul	1342140	Phenanthrene-d10	17.156

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	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048075.D
Acq On : 16 Oct 2024 00:22
Operator : RC/JU
Sample : P4350-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAJ9

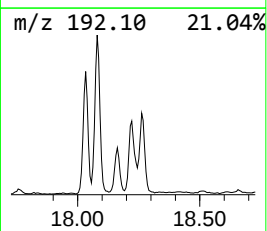
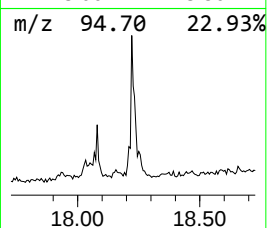
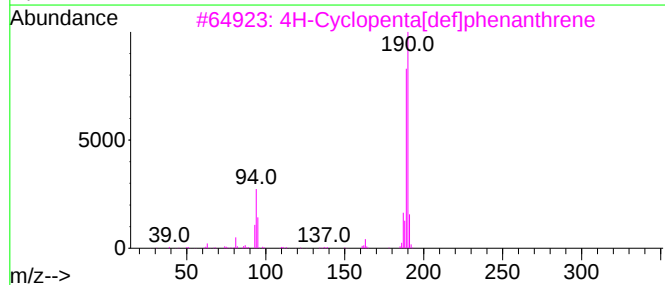
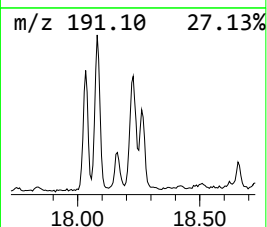
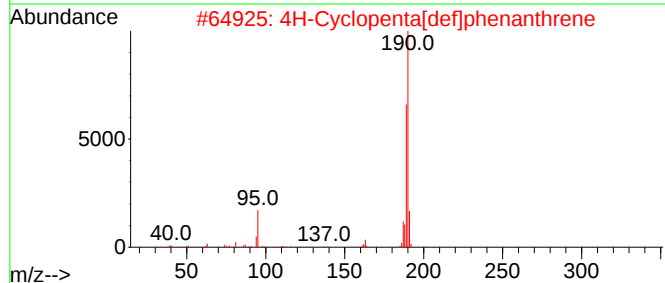
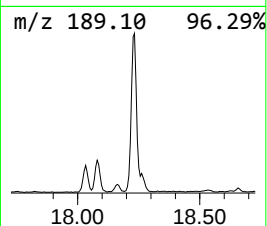
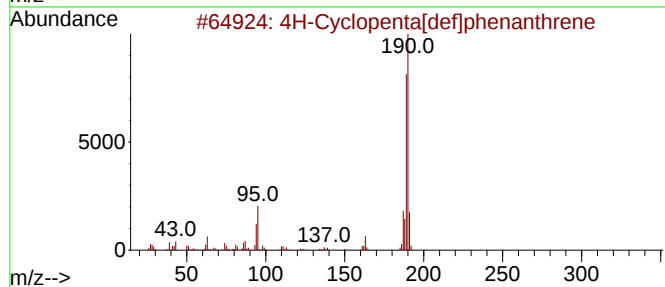
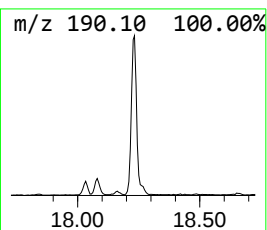
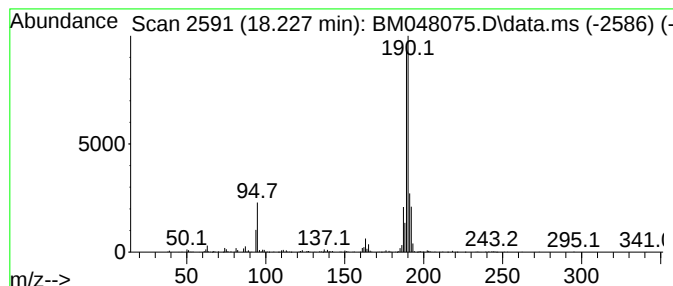
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	3.69 ng/ul	627753	Phenanthrene-d10	17.156

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	89
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
4			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	53
5			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048075.D
Acq On : 16 Oct 2024 00:22
Operator : RC/JU
Sample : P4350-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAJ9

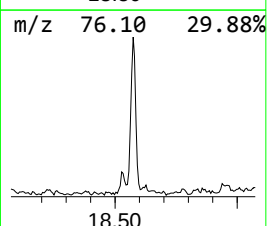
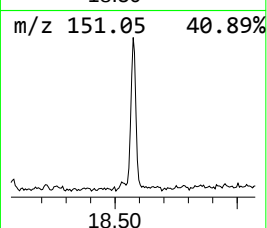
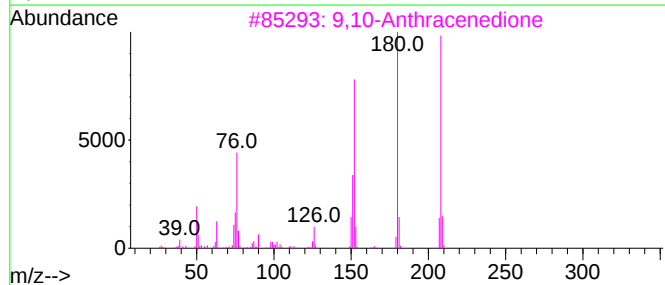
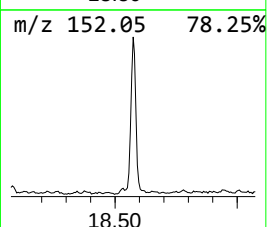
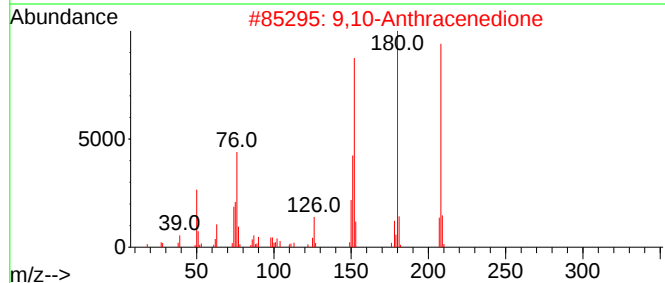
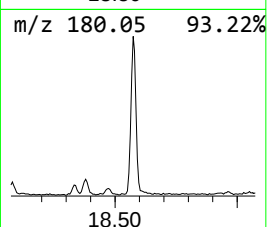
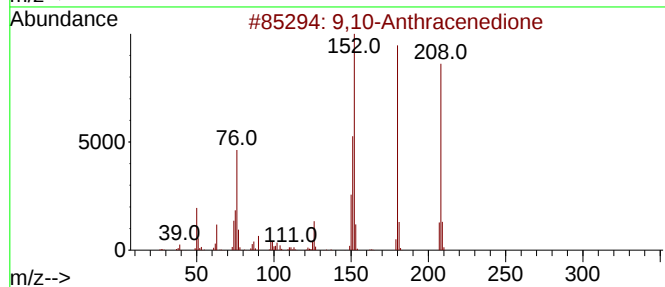
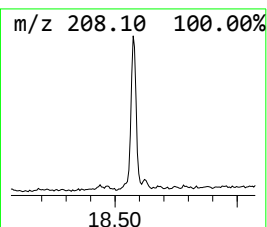
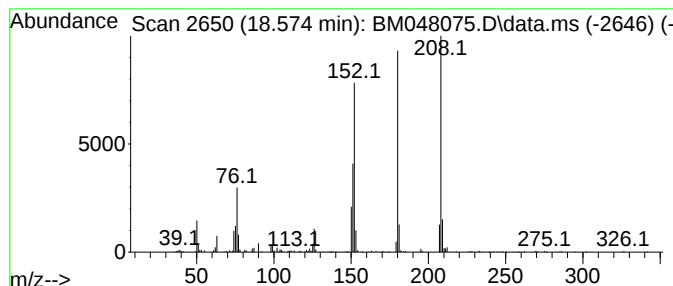
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.574	2.57 ng/ul	437136	Phenanthrene-d10	17.156

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	97
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	96
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048075.D
Acq On : 16 Oct 2024 00:22
Operator : RC/JU
Sample : P4350-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAJ9

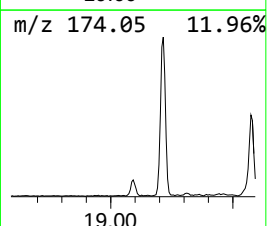
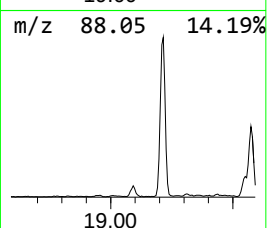
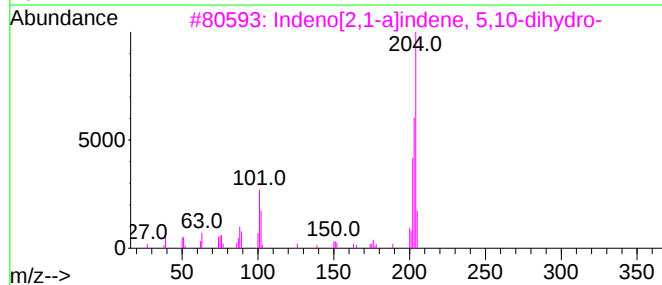
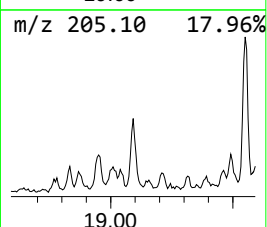
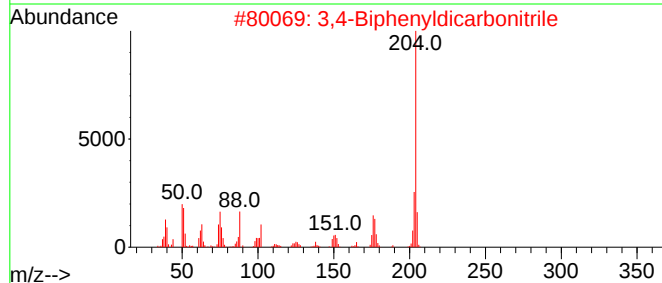
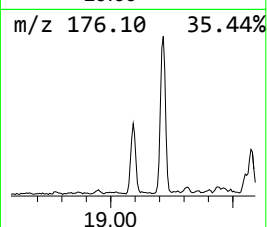
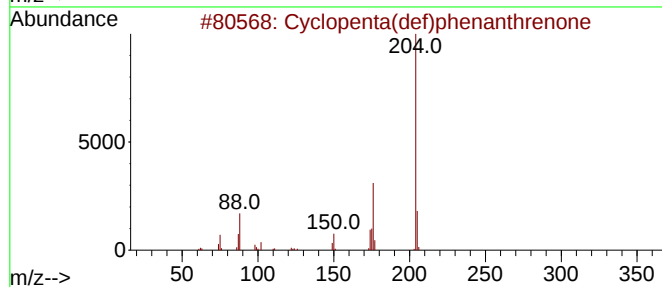
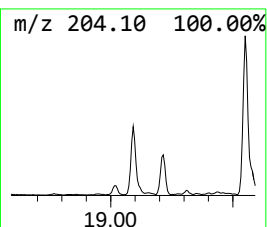
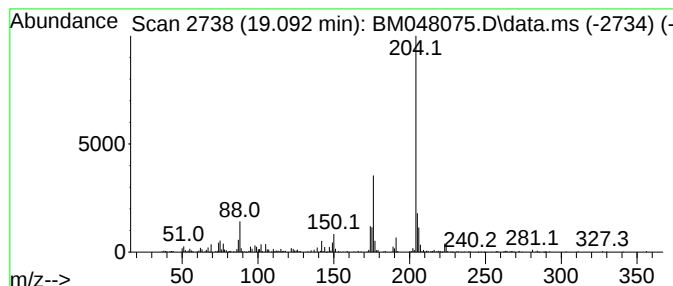
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.092	2.29 ng/ul	389359	Phenanthrene-d10	17.156

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	95
2			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	53
3			Indeno[2,1-a]indene, 5,10-dihydro-	204	C16H12	006543-29-9	50
4			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	50
5			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
Data File : BM048075.D
Acq On : 16 Oct 2024 00:22
Operator : RC/JU
Sample : P4350-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EOAJ9

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclobutane, et...	3.046	31.9	ng/ul	2630660	1	7.781	1646510	20.0
2-Cyclohexen-1-one	6.416	2.1	ng/ul	174016	1	7.781	1646510	20.0
Benzophenone	15.774	2.7	ng/ul	405148	3	14.415	2992400	20.0
n-Hexadecanoic ...	18.051	7.9	ng/ul	1342140	4	17.156	3401560	20.0
4H-Cyclopenta[d...	18.227	3.7	ng/ul	627753	4	17.156	3401560	20.0
9,10-Anthracene...	18.574	2.6	ng/ul	437136	4	17.156	3401560	20.0
Cyclopenta(def)...	19.092	2.3	ng/ul	389359	4	17.156	3401560	20.0