LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

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-BM121523.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM043372.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.387	30	34	39	rBV	124198	159229	1.00%	0.100%
2	3.669	80	82	93	rBV	106301	173593	1.09%	0.109%
3	3.805	102	105	123	rBV	1537621	2387809	14.98%	1.503%
4	5.481	385	390	397	rBV	56452	87871	0.55%	0.055%
5	5.610	408	412	421	rBV	52921	101128	0.63%	0.064%
6	7.151	666	674	681	rVV	2183892	3471487	21.77%	2.186%
7	7.334	698	705	720	rBV	1674862	2670913	16.75%	1.681%
8	7.522	730	737	747	rVV	2072869	3310781	20.77%	2.084%
9	7.634	751	756	763	rVV	71153	142556	0.89%	0.090%
10	7.710	763	769	779	rVB	50654	99019	0.62%	0.062%
11	7.993	809	817	825	rBV	1737634	2785664	17.47%	1.754%
12	8.363	872	880	890	rBV	1047058	1676642	10.52%	1.056%
13	8.528	901	908	917	rVB	97396	179361	1.12%	0.113%
14	8.704	931	938	949	rBV	2770169	4343687	27.24%	2.735%
15	8.851	956	963	970	rVB2	51178	88641	0.56%	0.056%
16	9.169	1009	1017	1026	rVV	2035214	3408833	21.38%	2.146%
17	9.351	1039	1048	1054	rVV2	37222	82599	0.52%	0.052%
18	9.481	1064	1070	1075	rVV	78494	169271	1.06%	0.107%
19	9.734	1108	1113	1121	rBV	73158	114163	0.72%	0.072%
20	9.886	1130	1139	1148	rBV	1694155	2777955	17.42%	1.749%
21	9.975	1148	1154	1157	rVV	112893	204523	1.28%	0.129%
22	10.010	1157	1160	1164	rVV	95174	156448	0.98%	0.098%
23	10.192	1183	1191	1201	rVV4	48265	158074	0.99%	0.100%
24	10.281	1201	1206	1223	rVV	303480	547720	3.44%	0.345%
25	10.422	1223	1230	1241	rVV	2575642	4369761	27.41%	2.751%
26	10.804	1286	1295	1299	rVV	2537892	4430010	27.79%	2.789%
27	10.857	1299	1304	1312	rVV	9614541	15943397	100.00%	10.037%
28	10.945	1312	1319	1335	rVB2	2514069	5126016	32.15%	3.227%
29	11.345	1379	1387	1392	rBV2	76236	149599	0.94%	0.094%
30	11.422	1392	1400	1406	rVB	62312	121927	0.76%	0.077%
31	11.551	1414	1422	1431	rBV	144906	269083	1.69%	0.169%
32	11.633	1431	1436	1444	rVV	187271	324426	2.03%	0.204%
33	11.828	1462	1469	1473	rBV	66282	126879	0.80%	0.080%
34	11.880	1473	1478	1482	rVV	76993	139636	0.88%	0.088%
35	12.192	1526	1531	1545	rVV4	63274	176413	1.11%	0.111%
36	12.298	1545	1549	1554	rVV3	41985	80562	0.51%	0.051%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

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-BM121523.MA.M
Title : SVOA CALIBRATION

37	12.351	1554	1558	1561	rVV	54318	97259	0.61%	0.061%
38	12.386	1561	1564	1569	rVV2	53199	98988	0.62%	0.062%
39	12.457	1569	1576	1586	rVV	1450316	2317861	14.54%	1.459%
40	12.563	1588	1594	1604	rVV7	98672	291401	1.83%	0.183%
41	12.675	1604	1613	1621	rVB	897396	1454950	9.13%	0.916%
42	12.845	1636	1642	1646	rBV4	46338	86353	0.54%	0.054%
43	13.463	1735	1747	1753	rBV2	266137	506662	3.18%	0.319%
44	13.522	1753	1757	1767	rVV	50744	105961	0.66%	0.067%
45	13.657	1776	1780	1790	rVB	39249	91275	0.57%	0.057%
46	13.804	1800	1805	1812	rVB3	59784	142961	0.90%	0.090%
47	13.945	1824	1829	1834	rVV	72819	129158	0.81%	0.081%
48	14.045	1840	1846	1854	rVV	4210644	5939943	37.26%	3.740%
49	14.322	1886	1893	1905	rVB	5038219	7595698	47.64%	4.782%
50	14.621	1936	1944	1950	rVV	3621125	5429702	34.06%	3.418%
51	14.686	1950	1955	1962	rVV	1268825	1812681	11.37%	1.141%
52	14.763	1962	1968	1972	rVV	141611	218439	1.37%	0.138%
53	14.821	1972	1978	1988	rVV	2096580	3156194	19.80%	1.987%
54	14.910	1988	1993	2001	rVV2	109387	215062	1.35%	0.135%
55	15.021	2007	2012	2021	rVV	401733	689889	4.33%	0.434%
56	15.616	2106	2113	2118	rBV	6012783	8604334	53.97%	5.417%
57	15.668	2118	2122	2127	rVB2	399883	542482	3.40%	0.342%
58	15.733	2127	2133	2141	rBV	1556890	2124477	13.33%	1.337%
59	15.974	2168	2174	2180	rBV2	135434	217287	1.36%	0.137%
60	16.110	2191	2197	2203	rVB4	44173	99833	0.63%	0.063%
61	16.345	2224	2237	2244	rBV2	297667	663001	4.16%	0.417%
62	16.533	2264	2269	2280	rVB4	62270	179208	1.12%	0.113%
63	16.633	2280	2286	2290	rBV2	68858	127473	0.80%	0.080%
64	16.768	2305	2309	2313	rBV2	66097	86490	0.54%	0.054%
65	17.015	2348	2351	2358	rVB3	57823	100143	0.63%	0.063%
66	17.180	2376	2379	2386	rVB2	66982	107254	0.67%	0.068%
67	17.368	2404	2411	2415	rVV	3687158	5500594	34.50%	3.463%
68	17.410	2415	2418	2422	rVV	594717	841245	5.28%	0.530%
69	17.462	2422	2427	2437	rVV2	5775518	8669940	54.38%	5.458%
70	17.557	2437	2443	2453	rVB4	118110	266223	1.67%	0.168%
71	17.751	2470	2476	2489	rBV2	1113344	2194017	13.76%	1.381%
72	18.268	2557	2564	2575	rBV2	1190592	1838365	11.53%	1.157%
73	18.439	2587	2593	2597	rBV2	151248	255295	1.60%	0.161%
74	18.692	2631	2636	2641	rBV	1688439	2014444	12.63%	1.268%
75	18.739	2641	2644	2648	rVV	66904	87606	0.55%	0.055%
76	18.780	2648	2651	2659	rVB	85651	115617	0.73%	0.073%

LSC Area Percent Report

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Sample : 05626-13
Misc :
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Instrument :
BNA_M
ClientSampleId :
DCLF9

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

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Title : SVOA CALIBRATION

77	18.980	2682	2685	2690	rVB	102308	128652	0.81%	0.081%
78	19.156	2712	2715	2719	rVB3	83894	106795	0.67%	0.067%
79	19.239	2726	2729	2734	rVB	206288	258269	1.62%	0.163%
80	19.415	2751	2759	2764	rBV	1309616	1855303	11.64%	1.168%
81	19.474	2765	2769	2773	rVB2	171341	203371	1.28%	0.128%
82	19.539	2776	2780	2789	rVB2	269321	423153	2.65%	0.266%
83	19.751	2810	2816	2820	rBV	6014978	8559188	53.68%	5.389%
84	20.021	2859	2862	2867	rVB	108650	143523	0.90%	0.090%
85	20.103	2873	2876	2881	rVB3	88801	153684	0.96%	0.097%
86	20.303	2907	2910	2917	rVB2	105939	120246	0.75%	0.076%
87	20.568	2952	2955	2958	rBV	102057	114599	0.72%	0.072%
88	20.798	2991	2994	2999	rVB2	124869	141621	0.89%	0.089%
89	21.015	3028	3031	3039	rVB	169698	223877	1.40%	0.141%
90	21.221	3063	3066	3068	rBV2	154575	178539	1.12%	0.112%
91	21.256	3069	3072	3078	rVV2	920522	1241818	7.79%	0.782%
92	21.539	3114	3120	3124	rBV	3268903	4967889	31.16%	3.128%
93	21.574	3124	3126	3133	rVB	767498	982088	6.16%	0.618%
94	21.709	3147	3149	3155	rVB2	243551	312366	1.96%	0.197%
95	22.180	3226	3229	3237	rVB4	282946	439487	2.76%	0.277%
96	22.827	3335	3339	3345	rVB	485948	723041	4.54%	0.455%
97	23.221	3401	3406	3411	rBV	920027	2083516	13.07%	1.312%
98	23.744	3491	3495	3499	rBV	429737	773381	4.85%	0.487%
99	23.803	3499	3505	3511	rVV2	2934958	5430155	34.06%	3.419%
100	23.962	3525	3532	3539	rVB	1865300	3705716	23.24%	2.333%

Sum of corrected areas: 158841718

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043372.D
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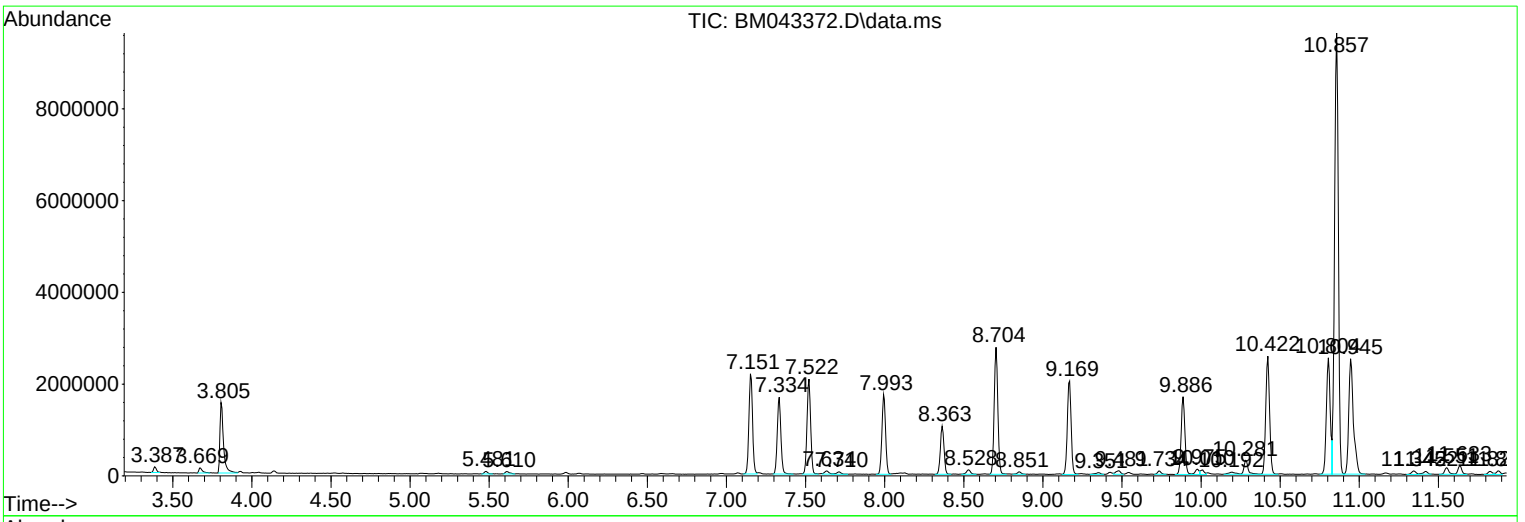
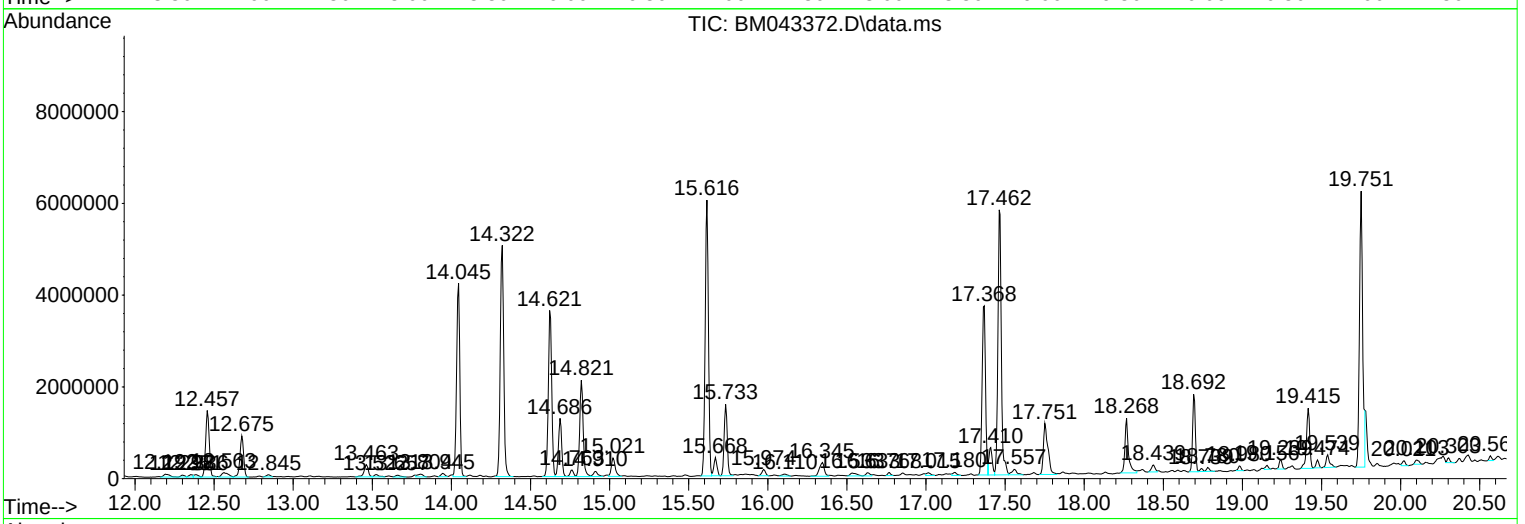
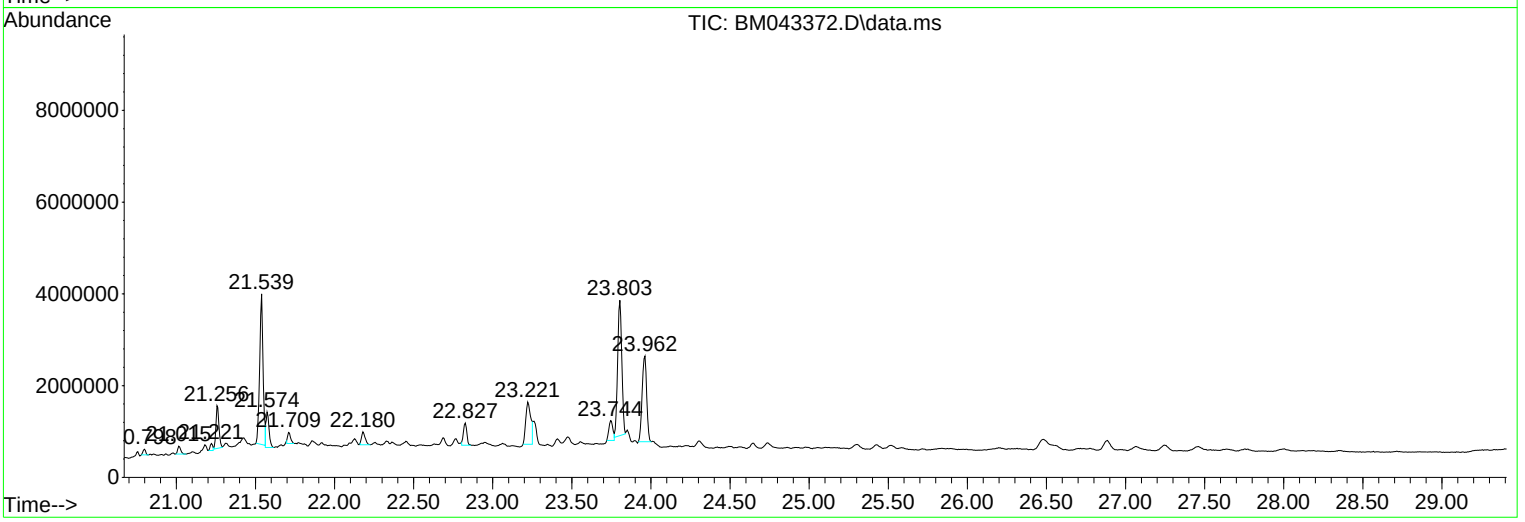
Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
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Sample : 05626-13
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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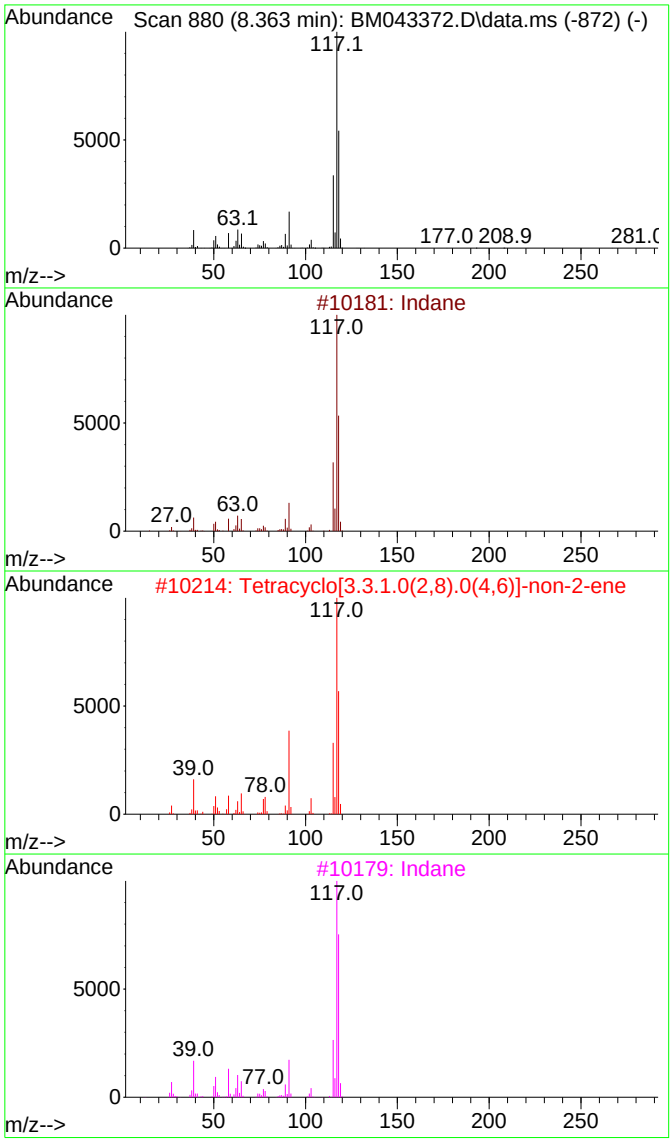
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

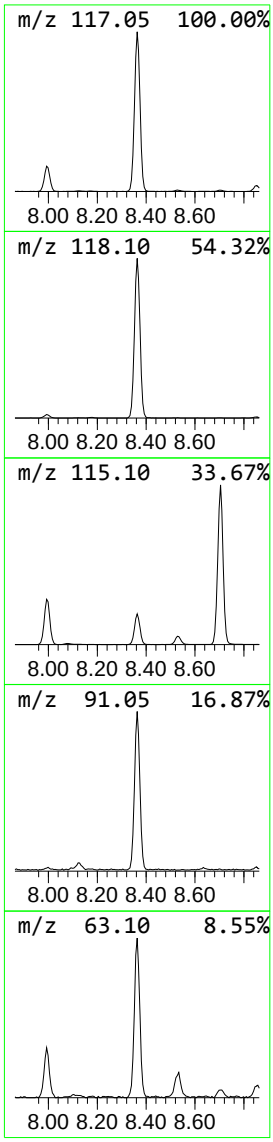
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.363	12.04 ng/ul	1676640	1,4-Dichlorobenzene-d4	7.992

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
3	Indane	118	C9H10	000496-11-7	64
4	Deltacyclene	118	C9H10	007785-10-6	64
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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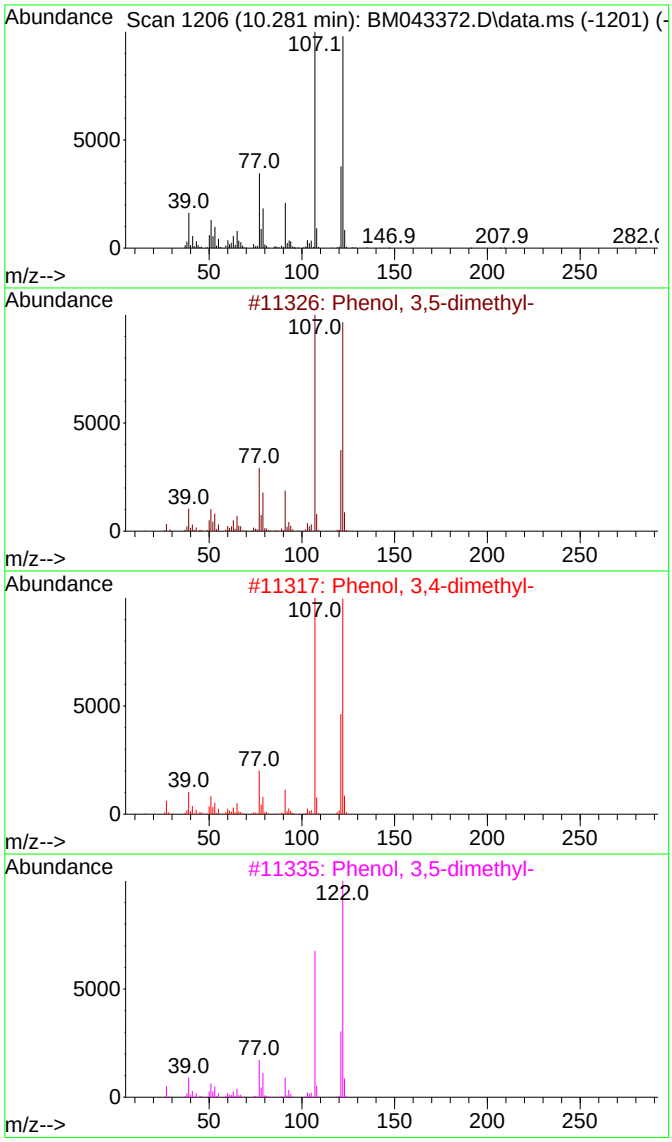
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

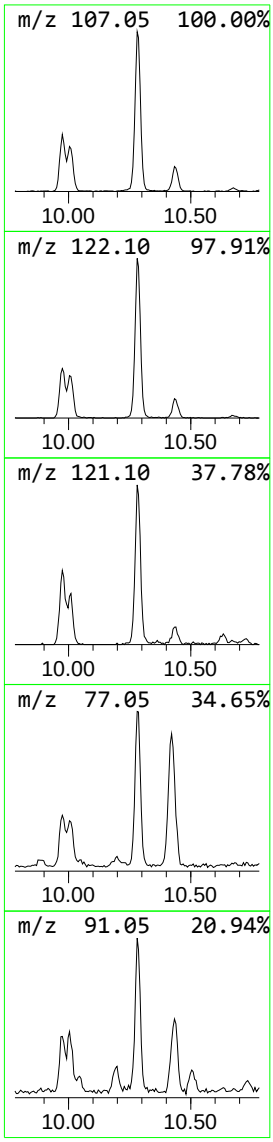
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Phenol, 3,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.281	2.47 ng/ul	547720	Naphthalene-d8	10.804

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94
5	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94





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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

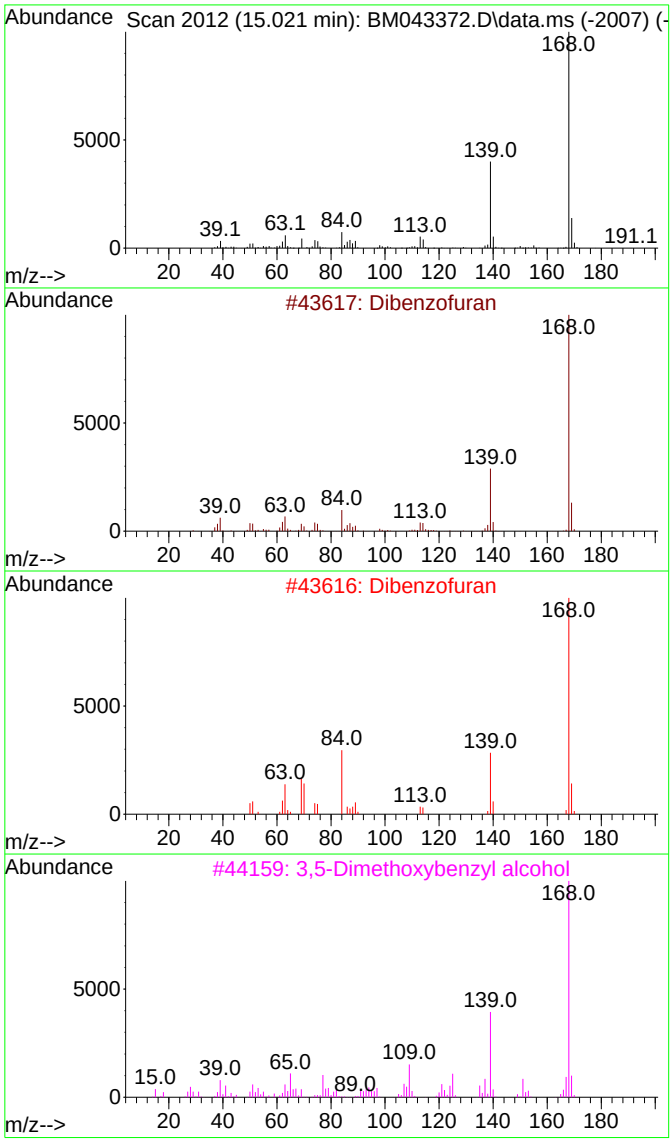
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

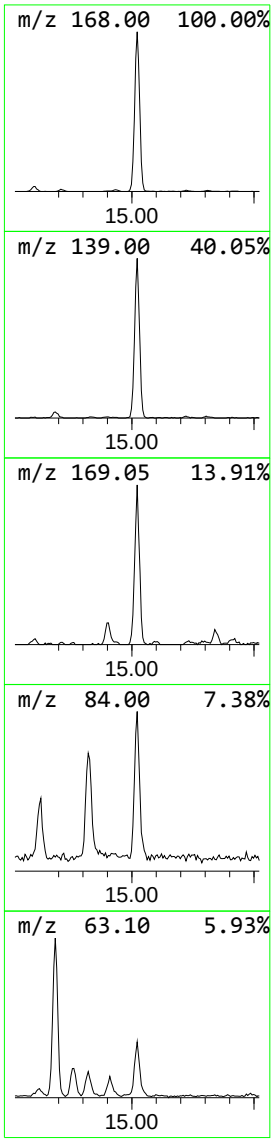
****

Peak Number 3 Dibenzo furan Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.022	2.54 ng/ul	689889	Acenaphthene-d10	14.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran	168	C12H8O	000132-64-9	91
2	Dibenzofuran	168	C12H8O	000132-64-9	78
3	3,5-Dimethoxybenzyl alcohol	168	C9H12O3	000705-76-0	56
4	Dibenzofuran	168	C12H8O	000132-64-9	53
5	Naphtho[2,1-d]imidazole	168	C11H8N2	000233-53-4	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
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Sample : 05626-13
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Instrument :
BNA_M
ClientSampleId :
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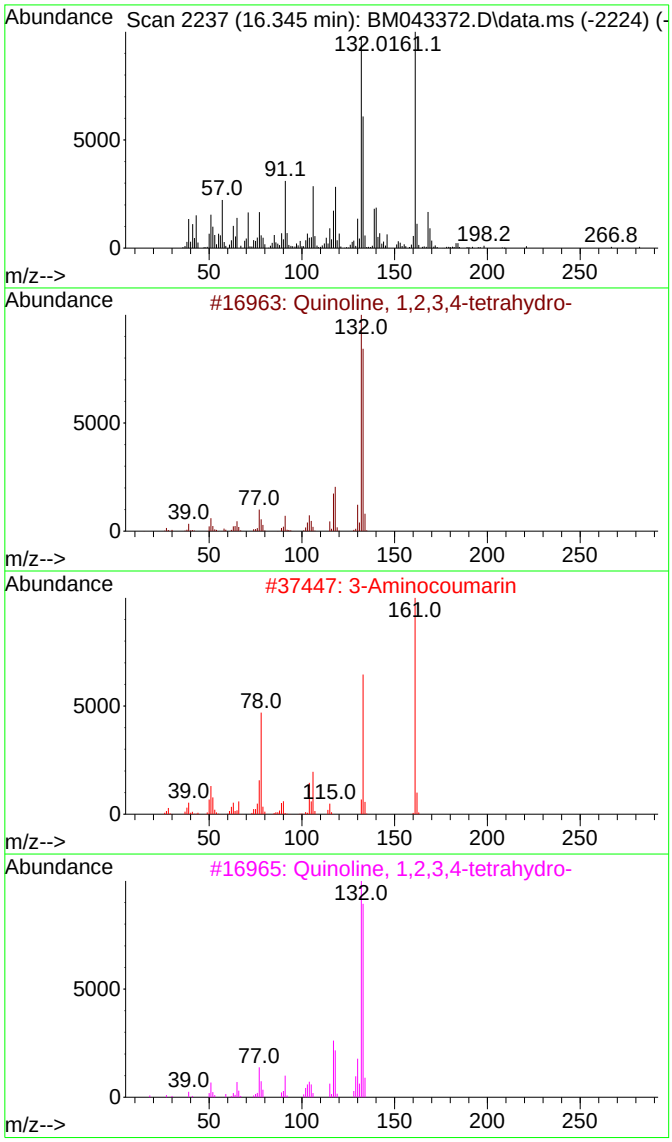
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

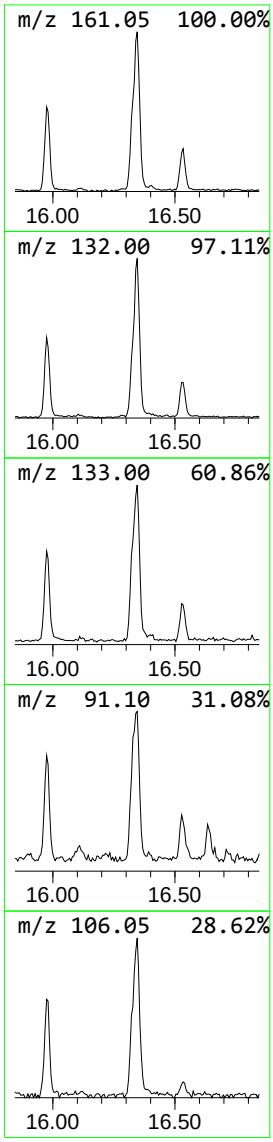
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Quinoline, 1,2,3,4-tetrahydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.345	2.41 ng/ul	663001	Phenanthrene-d10	17.368

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	64
2	3-Aminocoumarin	161	C9H7NO2	001635-31-0	50
3	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	46
4	5-Aminoindan	133	C9H11N	024425-40-9	46
5	Isoquinoline, 5,6,7,8-tetrahydro-	133	C9H11N	036556-06-6	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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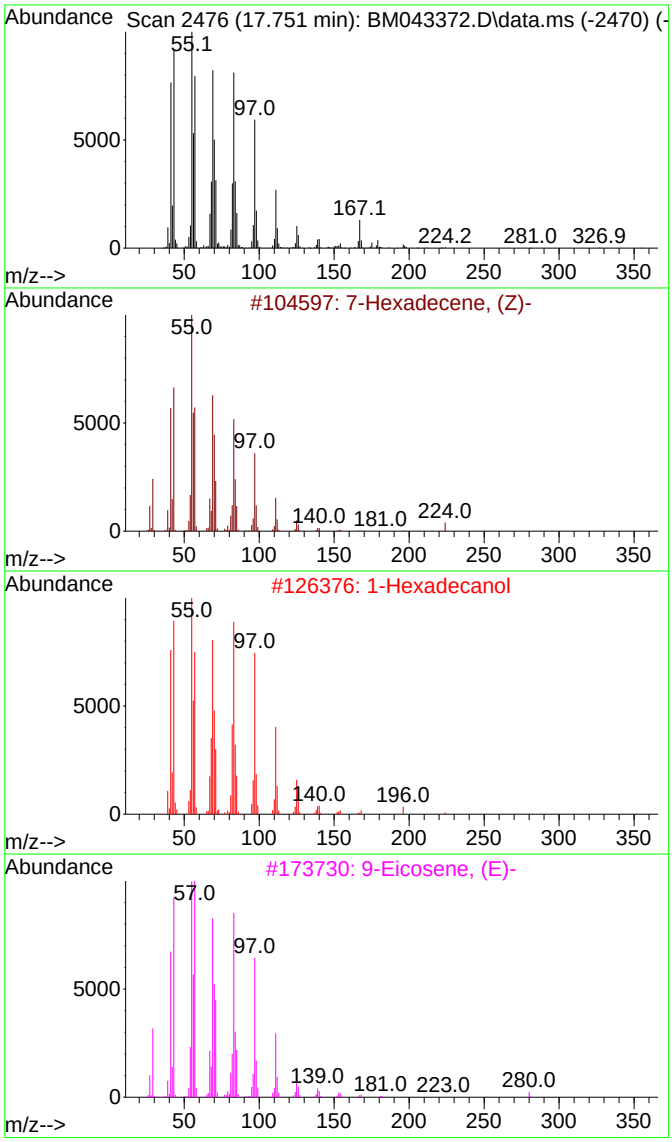
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

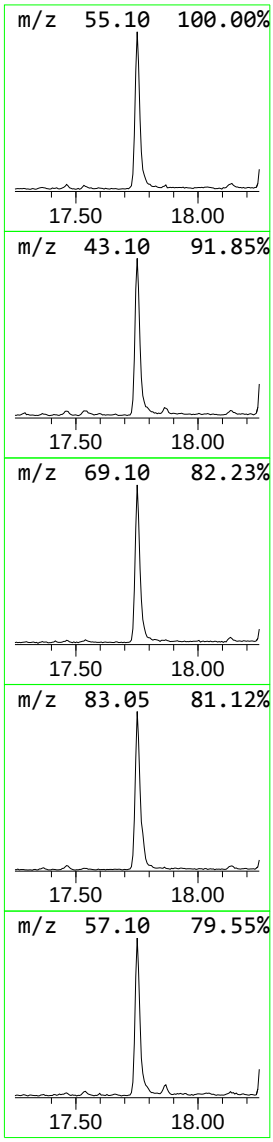
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.751	7.98 ng/ul	2194020	Phenanthrene-d10	17.368

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Hexadecene, (Z)-	224	C16H32	035507-09-6	98
2	1-Hexadecanol	242	C16H34O	036653-82-4	94
3	9-Eicosene, (E)-	280	C20H40	074685-29-3	94
4	1-Hexadecanol	242	C16H34O	036653-82-4	94
5	Cyclotetradecane	196	C14H28	000295-17-0	93





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Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

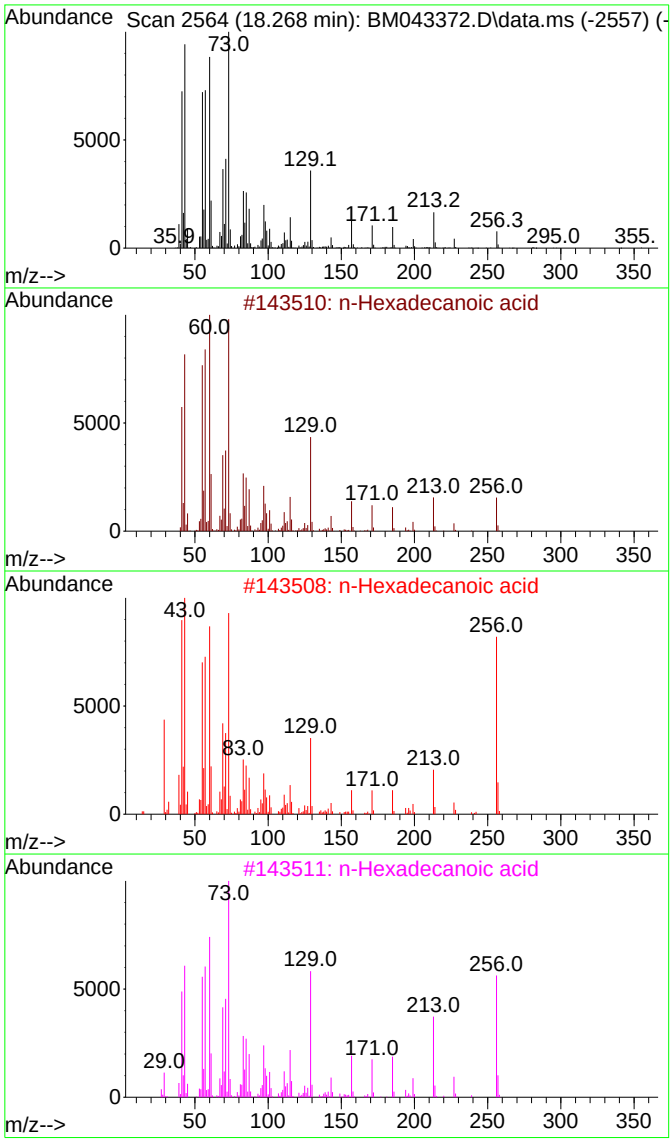
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

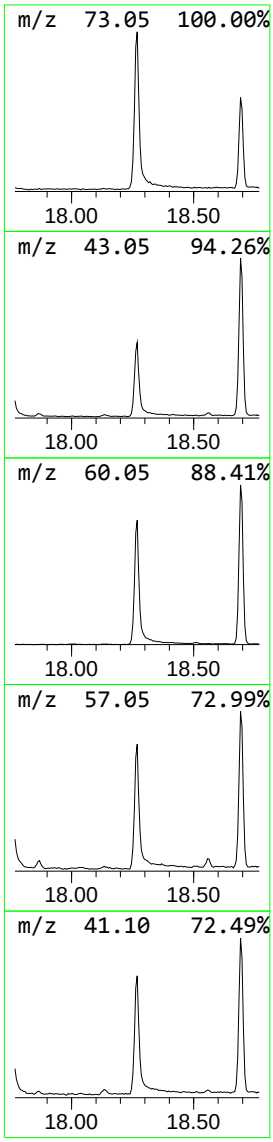
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.268	6.68 ng/ul	1838370	Phenanthrene-d10	17.368

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	99
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	94





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Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

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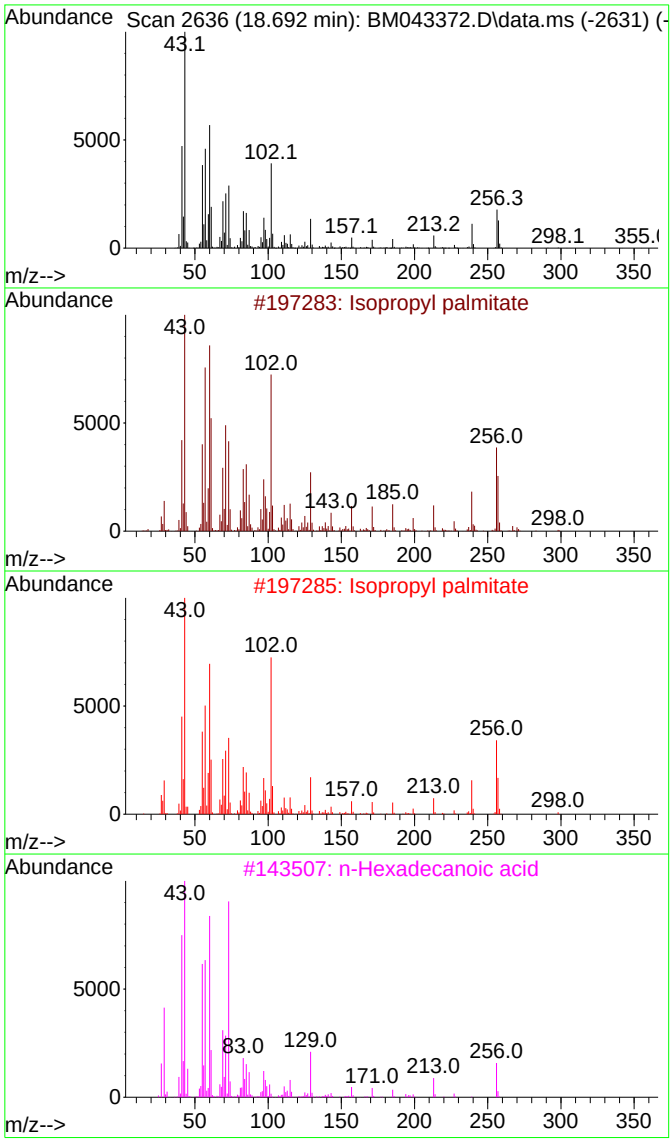
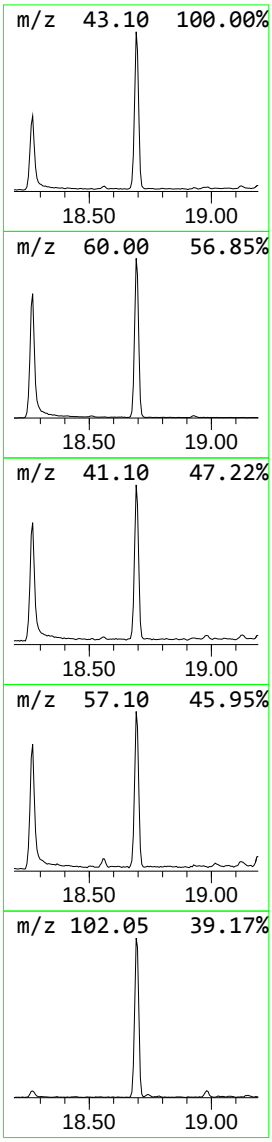
Peak Number 7 Isopropyl palmitate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.692	7.32 ng/ul	2014440	Phenanthrene-d10	17.368

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl palmitate	298	C19H38O2	000142-91-6	96
2	Isopropyl palmitate	298	C19H38O2	000142-91-6	91
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	49
4	n-Capric acid isopropyl ester	214	C13H26O2	002311-59-3	43
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	38





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Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

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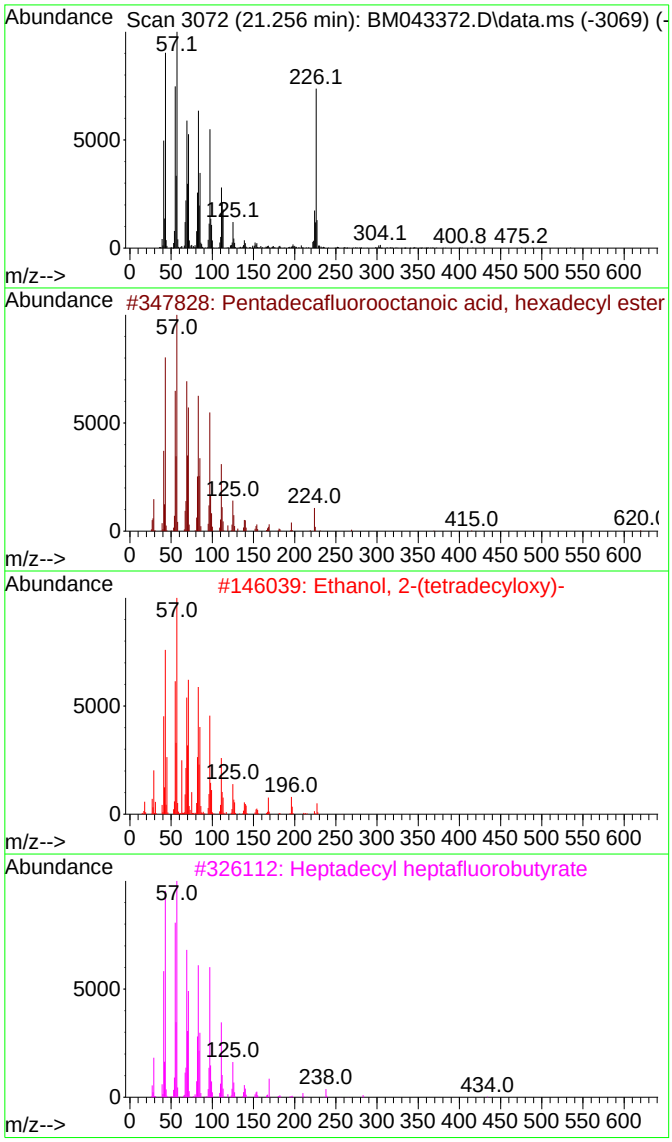
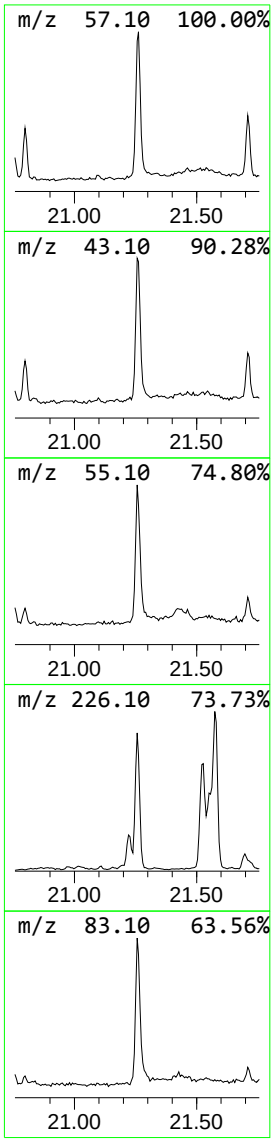
Peak Number 8 Pentadecafluorooctanoic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.256	5.00 ng/ul	1241820	Chrysene-d12	21.539

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	78
2	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	60
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	60
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	60
5	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	60





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Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

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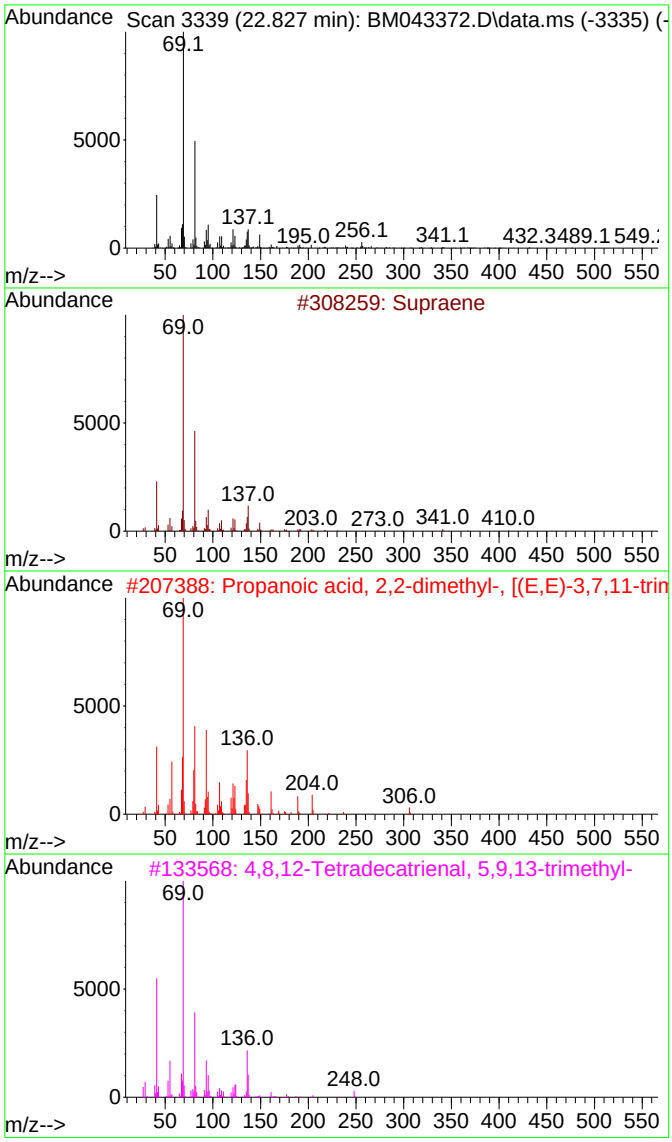
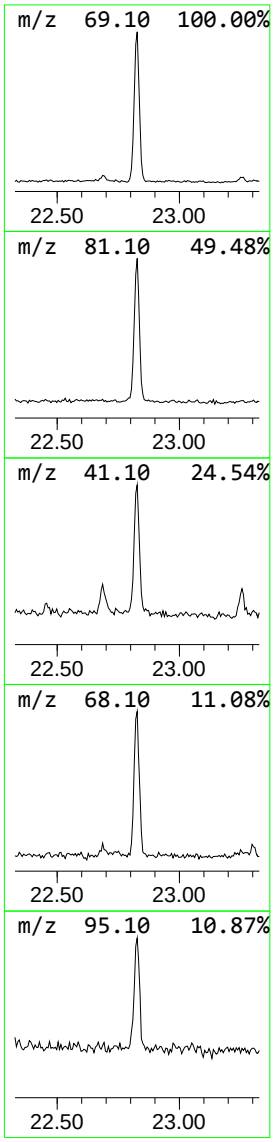
Peak Number 9 Supraene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.827	3.90 ng/ul	723041	Perylene-d12	23.962

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	91
2	Propanoic acid, 2,2-dimethyl-, [...	306	C20H34O2	1000164-38-8	78
3	4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
4	4,8,12-Tetradecatrienenitrile, 5...	245	C17H27N	005981-31-7	64
5	1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	58





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Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
Sample : 05626-13
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

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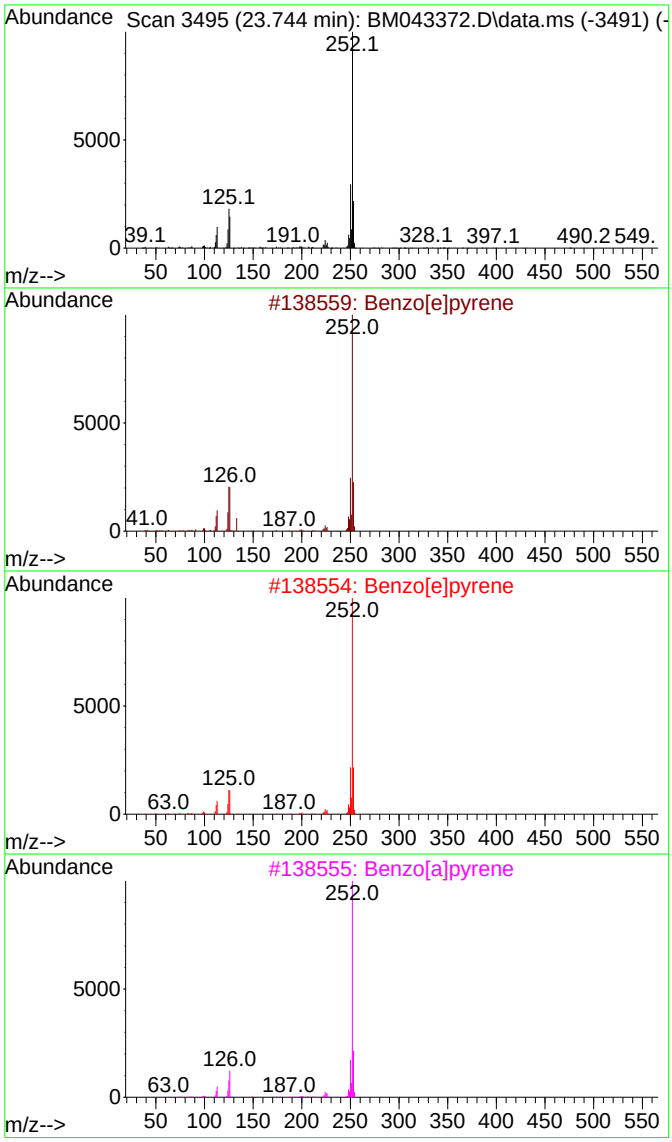
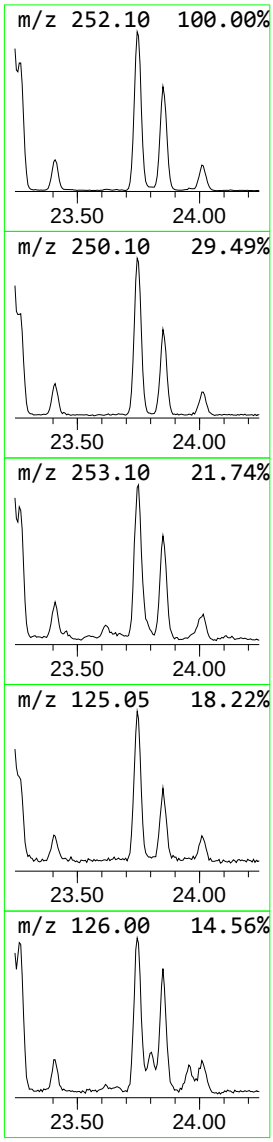
Peak Number 10 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.744	4.17 ng/ul	773381	Perylene-d12	23.962

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[e]pyrene	252	C20H12	000192-97-2	98
3	Benzo[a]pyrene	252	C20H12	000050-32-8	96
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	96
5	Perylene	252	C20H12	000198-55-0	94





Tentatively Identified Compound (LSC) summary

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Data File : BM043372.D
Acq On : 15 Dec 2023 23:49
Operator : MA/JU
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Misc :
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Instrument :
BNA_M
ClientSampleId :
DCLF9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	8.363	12.0	ng/ul	1676640	1	7.992	2785660	20.0
Phenol, 3,5-dim...	10.281	2.5	ng/ul	547720	2	10.804	4430010	20.0
Dibenzo furan	15.022	2.5	ng/ul	689889	3	14.621	5429700	20.0
Quinoline, 1,2,...	16.345	2.4	ng/ul	663001	4	17.368	5500590	20.0
(DEL) Alkane: S...	17.751	8.0	ng/ul	2194020	4	17.368	5500590	20.0
n-Hexadecanoic ...	18.268	6.7	ng/ul	1838370	4	17.368	5500590	20.0
Isopropyl palmi...	18.692	7.3	ng/ul	2014440	4	17.368	5500590	20.0
Pentadecafluoro...	21.256	5.0	ng/ul	1241820	5	21.539	4967890	20.0
Supraene	22.827	3.9	ng/ul	723041	6	23.962	3705720	20.0
Benzo[e]pyrene	23.744	4.2	ng/ul	773381	6	23.962	3705720	20.0