LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

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_G\METHODS\SFAM-EPA-BG090721.M
Title : SVOA CALIBRATION

Signal : TIC: BG050239.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.066	171	175	183	rVB	21942	34460	1.02%	0.211%
2	5.417	400	405	411	rBV	74441	119489	3.53%	0.733%
3	5.552	423	428	438	rVB	36230	83953	2.48%	0.515%
4	5.928	486	492	502	rVB	46284	83708	2.47%	0.513%
5	7.267	715	720	726	rBV2	11184	18967	0.56%	0.116%
6	7.485	752	757	763	rVB3	11073	17872	0.53%	0.110%
7	7.584	770	774	781	rVB2	23985	37354	1.10%	0.229%
8	7.661	781	787	797	rBV	60583	105623	3.12%	0.648%
9	7.949	829	836	845	rVB	106973	182339	5.38%	1.118%
10	8.060	845	855	861	rBV3	8567	17136	0.51%	0.105%
11	8.319	891	899	907	rBV	94703	157896	4.66%	0.968%
12	8.483	919	927	937	rVB	344540	589468	17.40%	3.615%
13	8.683	956	961	968	rVV3	14858	24776	0.73%	0.152%
14	8.806	978	982	987	rVB2	10397	16057	0.47%	0.098%
15	9.123	1031	1036	1043	rBV4	19350	35917	1.06%	0.220%
16	9.311	1062	1068	1075	rBV2	16106	25405	0.75%	0.156%
17	9.435	1079	1089	1095	rBV2	36660	78894	2.33%	0.484%
18	9.852	1155	1160	1166	rBV3	12096	23164	0.68%	0.142%
19	9.999	1180	1185	1196	rVB4	13355	26068	0.77%	0.160%
20	10.157	1206	1212	1220	rBV4	25089	52665	1.55%	0.323%
21	10.392	1246	1252	1262	rBV3	44863	99519	2.94%	0.610%
22	10.692	1296	1303	1308	rBV8	12631	34671	1.02%	0.213%
23	10.762	1308	1315	1318	rVV	246590	437996	12.93%	2.686%
24	10.815	1318	1324	1336	rVB	1817067	3387604	100.00%	20.774%
25	10.939	1336	1345	1351	rBV2	163412	348604	10.29%	2.138%
26	12.172	1548	1555	1562	rVB6	14556	32715	0.97%	0.201%
27	12.319	1575	1580	1587	rVB2	25344	43349	1.28%	0.266%
28	12.431	1593	1599	1604	rVB	42927	74713	2.21%	0.458%
29	12.548	1614	1619	1623	rVV3	11590	20227	0.60%	0.124%
30	12.642	1625	1635	1643	rVB	309517	551095	16.27%	3.379%
31	12.842	1665	1669	1676	rVB6	13299	21337	0.63%	0.131%
32	13.071	1703	1708	1717	rVB8	15128	30921	0.91%	0.190%
33	13.353	1747	1756	1762	rBV3	21603	41950	1.24%	0.257%
34	13.435	1762	1770	1780	rVB2	88288	170816	5.04%	1.047%
35	13.576	1787	1794	1799	rBV3	28460	47330	1.40%	0.290%
36	13.635	1799	1804	1815	rVB5	24568	52421	1.55%	0.321%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

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_G\METHODS\SFAM-EPA-BG090721.M
Title : SVOA CALIBRATION

37	13.782	1818	1829	1836	rBV3	84842	177046	5.23%	1.086%
38	13.923	1848	1853	1858	rBV4	39816	75940	2.24%	0.466%
39	13.987	1858	1864	1865	rVV2	27835	47415	1.40%	0.291%
40	14.005	1865	1867	1874	rVB	45976	63694	1.88%	0.391%
41	14.164	1888	1894	1898	rBV3	16583	28577	0.84%	0.175%
42	14.299	1910	1917	1928	rBV2	48579	122376	3.61%	0.750%
43	14.604	1961	1969	1975	rVV	257634	439200	12.96%	2.693%
44	14.669	1975	1980	1987	rVB	348765	557411	16.45%	3.418%
45	14.745	1987	1993	1999	rBV	71453	109840	3.24%	0.674%
46	14.804	1999	2003	2010	rVB3	16809	29238	0.86%	0.179%
47	14.880	2011	2016	2020	rBV	27634	40305	1.19%	0.247%
48	15.004	2031	2037	2046	rVV	178889	290587	8.58%	1.782%
49	15.162	2057	2064	2072	rBV2	311987	530180	15.65%	3.251%
50	15.245	2073	2078	2087	rVV2	85155	171573	5.06%	1.052%
51	15.327	2089	2092	2100	rVV5	9167	19878	0.59%	0.122%
52	15.603	2133	2139	2143	rBV	65836	93030	2.75%	0.570%
53	15.656	2143	2148	2155	rVB	181917	260297	7.68%	1.596%
54	15.767	2159	2167	2172	rBV3	78989	174600	5.15%	1.071%
55	15.891	2181	2188	2195	rVB9	18979	57301	1.69%	0.351%
56	15.950	2195	2198	2203	rVB4	11634	16119	0.48%	0.099%
57	16.091	2218	2222	2231	rVB7	15384	35827	1.06%	0.220%
58	16.343	2258	2265	2270	rBV3	21739	39178	1.16%	0.240%
59	16.549	2295	2300	2302	rVV3	21909	40125	1.18%	0.246%
60	16.584	2302	2306	2310	rVV	67615	117800	3.48%	0.722%
61	16.619	2310	2312	2317	rVV	32796	50180	1.48%	0.308%
62	16.660	2317	2319	2325	rVV5	10287	17645	0.52%	0.108%
63	16.713	2325	2328	2334	rVB5	12525	24155	0.71%	0.148%
64	16.872	2347	2355	2363	rVB5	26389	53087	1.57%	0.326%
65	17.019	2372	2380	2391	rBV	650187	1139434	33.64%	6.987%
66	17.148	2399	2402	2405	rBV2	20002	24819	0.73%	0.152%
67	17.177	2405	2407	2415	rVB	22165	29025	0.86%	0.178%
68	17.254	2415	2420	2424	rBV3	11504	21634	0.64%	0.133%
69	17.359	2433	2438	2442	rBV	350645	519634	15.34%	3.187%
70	17.406	2442	2446	2451	rVV	140876	226094	6.67%	1.386%
71	17.459	2451	2455	2460	rVB	68115	96631	2.85%	0.593%
72	17.565	2468	2473	2477	rBV6	13718	18079	0.53%	0.111%
73	17.694	2489	2495	2498	rBV3	12266	16389	0.48%	0.101%
74	17.771	2498	2508	2514	rVB	487250	761210	22.47%	4.668%
75	17.865	2519	2524	2531	rVV4	38693	80472	2.38%	0.493%
76	17.929	2531	2535	2545	rVV	56135	105393	3.11%	0.646%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

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_G\METHODS\SFAM-EPA-BG090721.M
Title : SVOA CALIBRATION

77	18.011	2545	2549	2555	rVV4	20977	40594	1.20%	0.249%
78	18.070	2555	2559	2563	rVV6	13469	22390	0.66%	0.137%
79	18.123	2563	2568	2574	rVB4	11208	20156	0.59%	0.124%
80	18.205	2574	2582	2590	rVB	33568	58656	1.73%	0.360%
81	18.288	2591	2596	2605	rBV2	49843	86243	2.55%	0.529%
82	18.364	2605	2609	2616	rVV	36863	57374	1.69%	0.352%
83	18.440	2617	2622	2627	rVB4	16548	25652	0.76%	0.157%
84	18.534	2631	2638	2644	rBV2	25442	68998	2.04%	0.423%
85	18.587	2644	2647	2651	rVV	16348	23952	0.71%	0.147%
86	18.675	2658	2662	2669	rVV2	18419	36113	1.07%	0.221%
87	18.740	2669	2673	2677	rVV2	12910	16489	0.49%	0.101%
88	19.357	2773	2778	2784	rBV2	19681	29914	0.88%	0.183%
89	19.415	2784	2788	2795	rVB3	10562	20258	0.60%	0.124%
90	19.750	2840	2845	2850	rBV	95544	139208	4.11%	0.854%
91	20.161	2910	2915	2920	rBV	22778	35067	1.04%	0.215%
92	20.226	2921	2926	2933	rVV	81089	117912	3.48%	0.723%
93	20.878	3034	3037	3042	rVB3	17891	28749	0.85%	0.176%
94	21.630	3159	3165	3171	rBV	397983	624684	18.44%	3.831%
95	23.874	3544	3547	3554	rVB4	13904	29198	0.86%	0.179%
96	24.620	3666	3674	3685	rVB3	50809	141907	4.19%	0.870%
97	24.838	3701	3711	3725	rVB	233626	692556	20.44%	4.247%
98	25.425	3808	3811	3813	rBV4	11834	16612	0.49%	0.102%
99	25.936	3892	3898	3908	rVB4	36411	106840	3.15%	0.655%
100	26.565	4002	4005	4008	rBV5	10717	19637	0.58%	0.120%

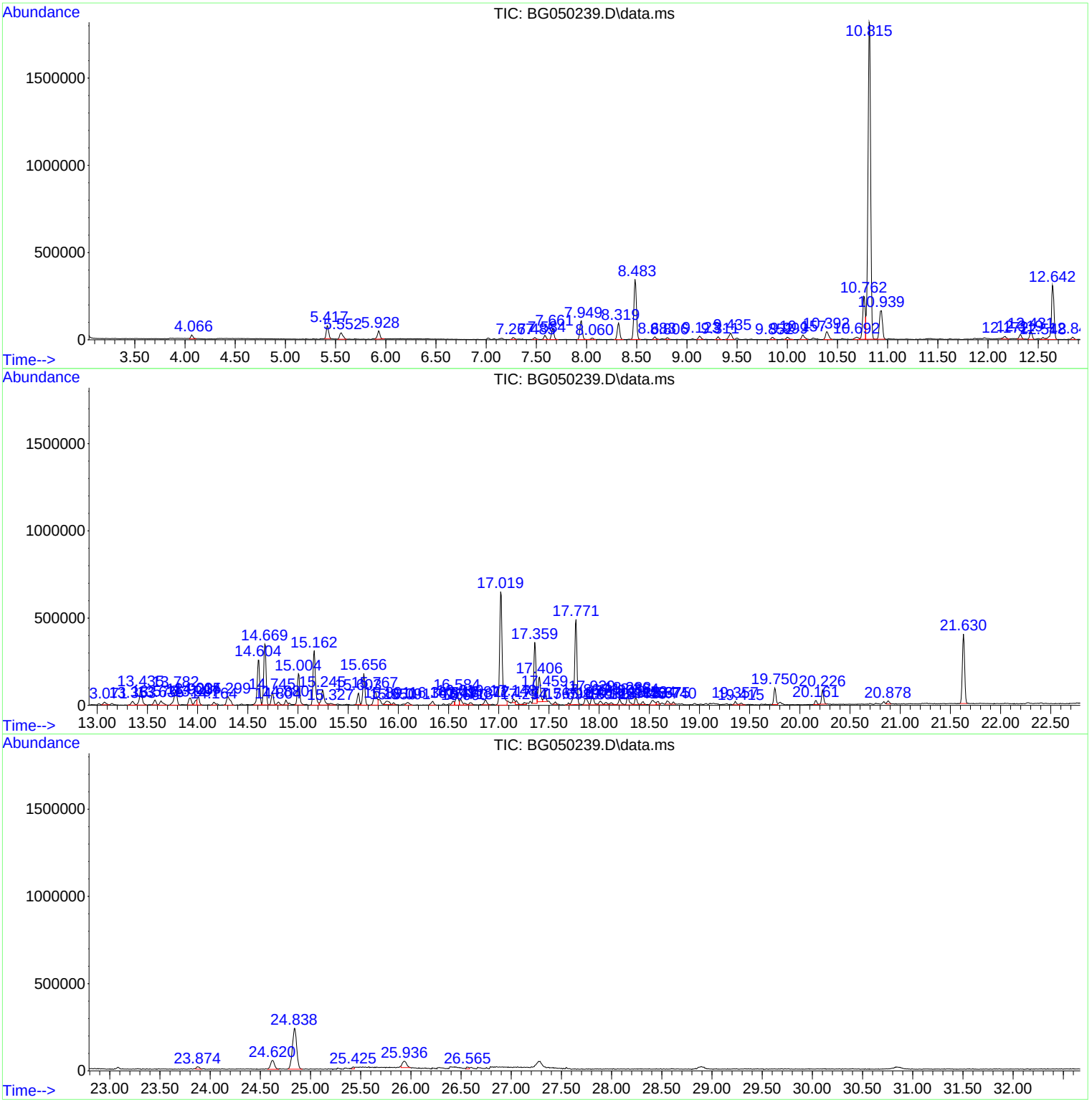
Sum of corrected areas: 16307056

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

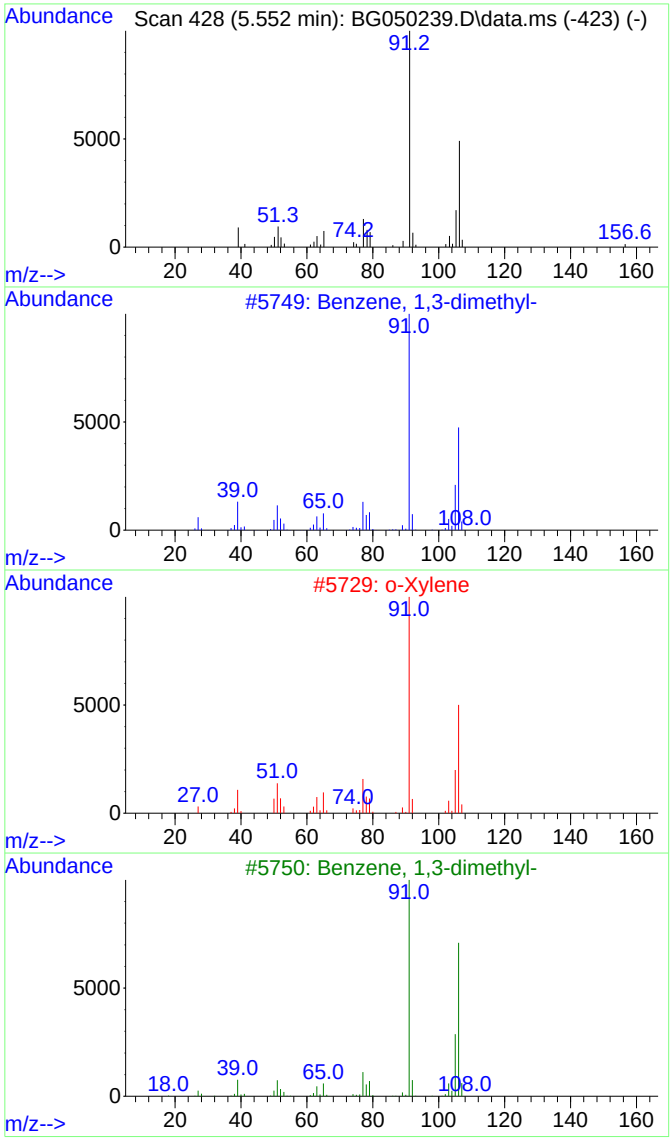
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

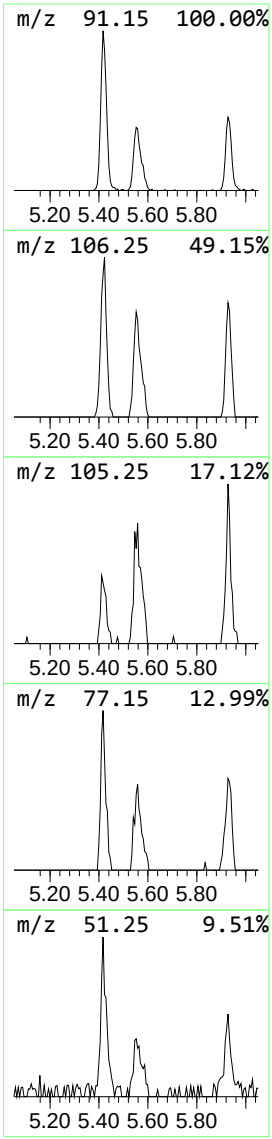
****

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.552	9.21 ng/ul	83953	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	o-Xylene	106	C8H10	000095-47-6	95
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
4	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
5	p-Xylene	106	C8H10	000106-42-3	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

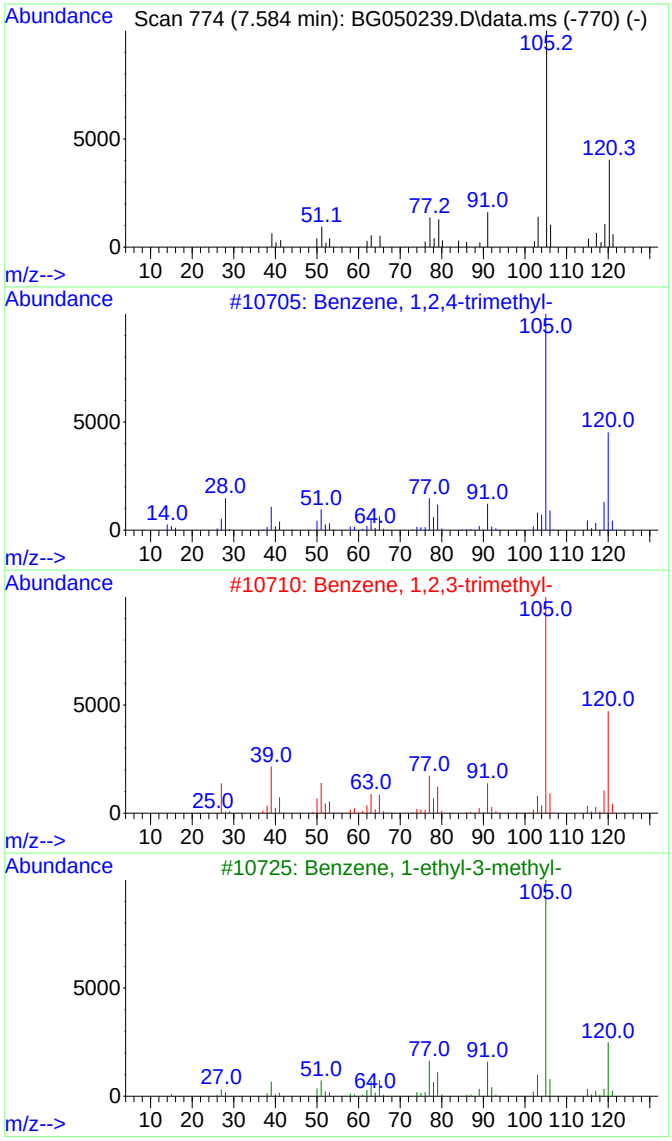
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

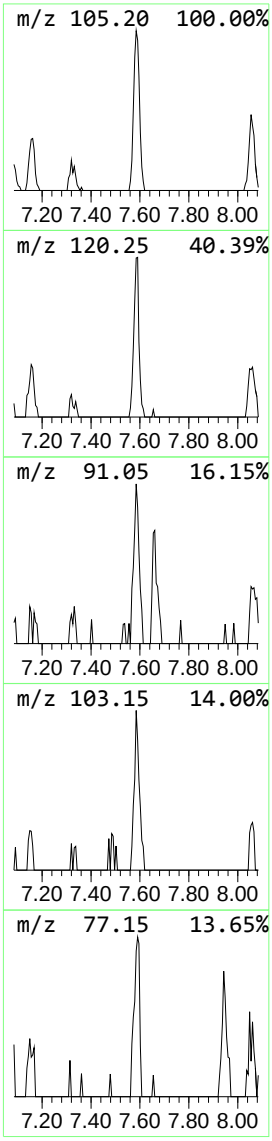
****

Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.584	4.10 ng/ul	37354	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

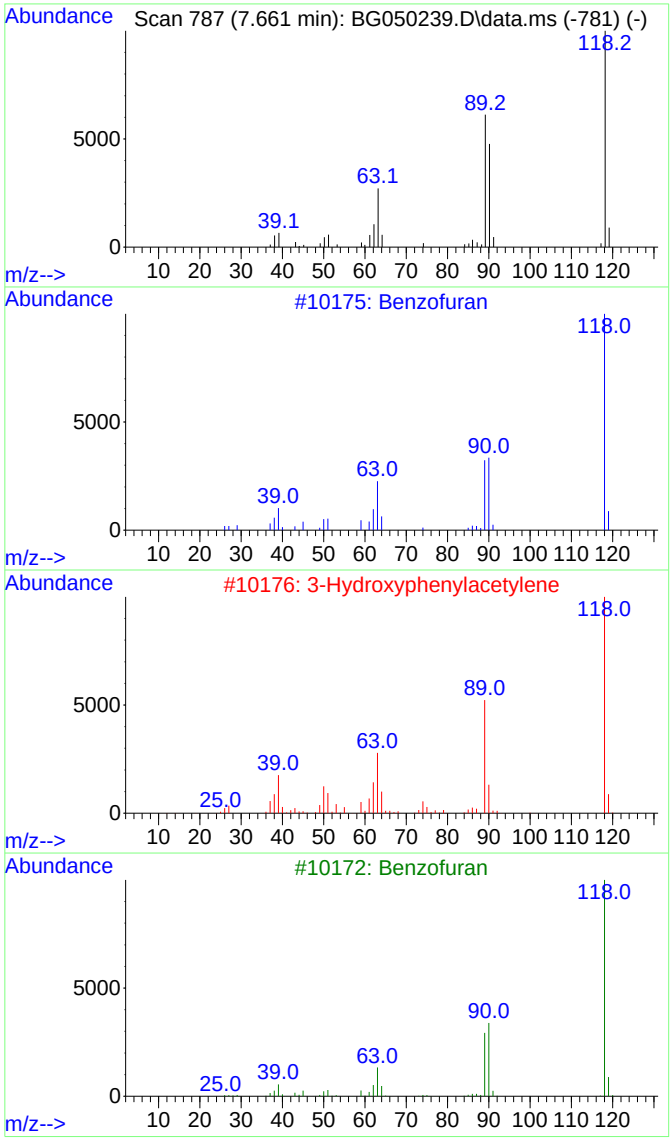
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

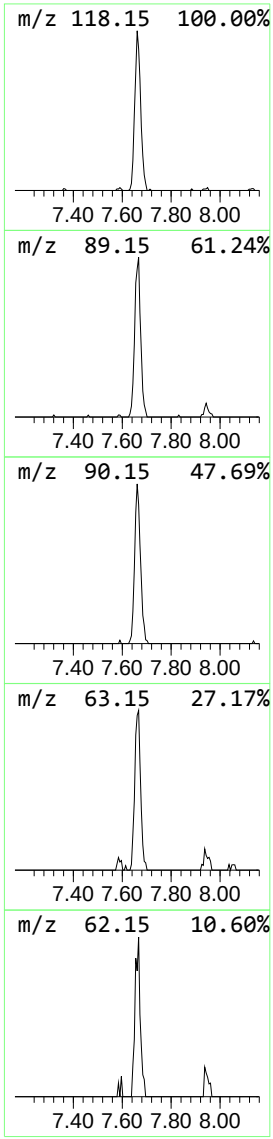
****

Peak Number 6 Benzofuran Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.661	11.59 ng/ul	105623	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	81
2	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	80
3	Benzofuran	118	C8H6O	000271-89-6	72
4	Benzofuran	118	C8H6O	000271-89-6	72
5	2H-Cyclopenta[d]pyridazine	118	C7H6N2	000270-64-4	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

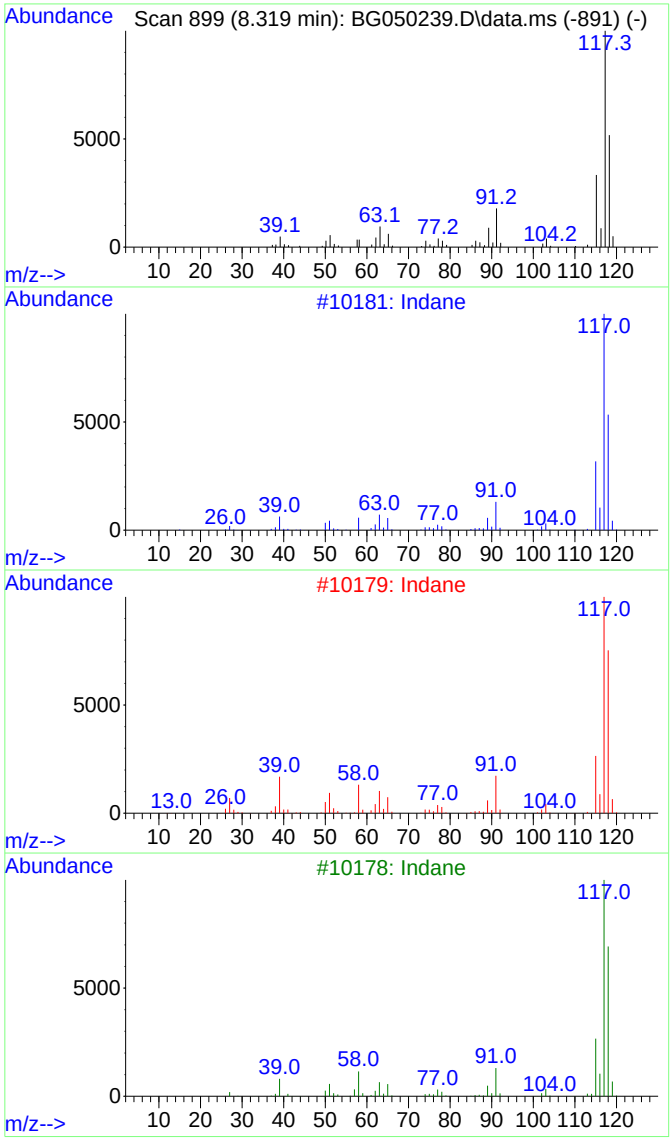
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

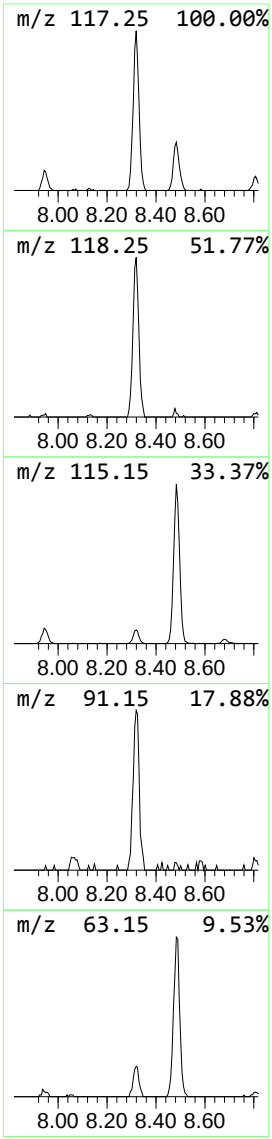
****

Peak Number 7 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.319	17.32 ng/ul	157896	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	81
4	Deltacyclene	118	C9H10	007785-10-6	68
5	Indane	118	C9H10	000496-11-7	60





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

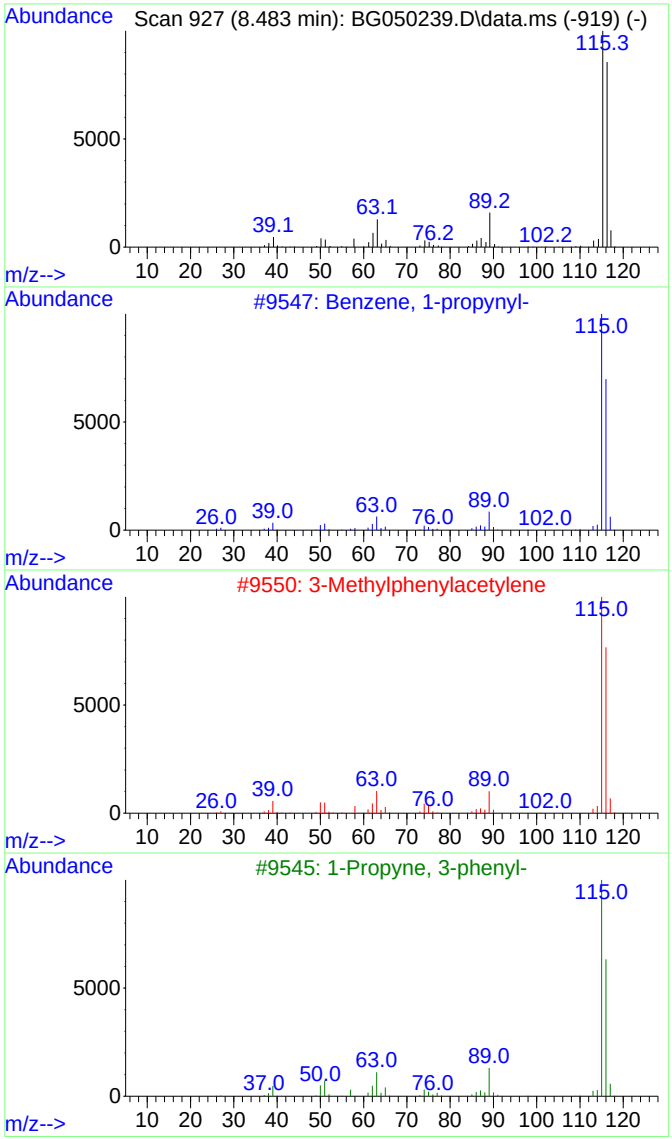
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

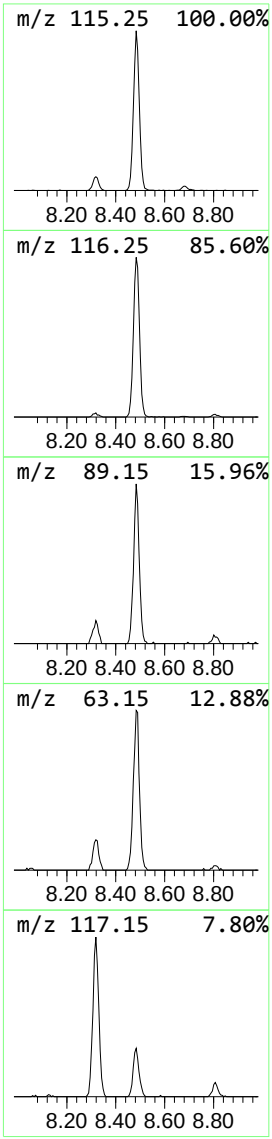
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-propynyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.483	64.66 ng/ul	589468	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
2	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
3	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

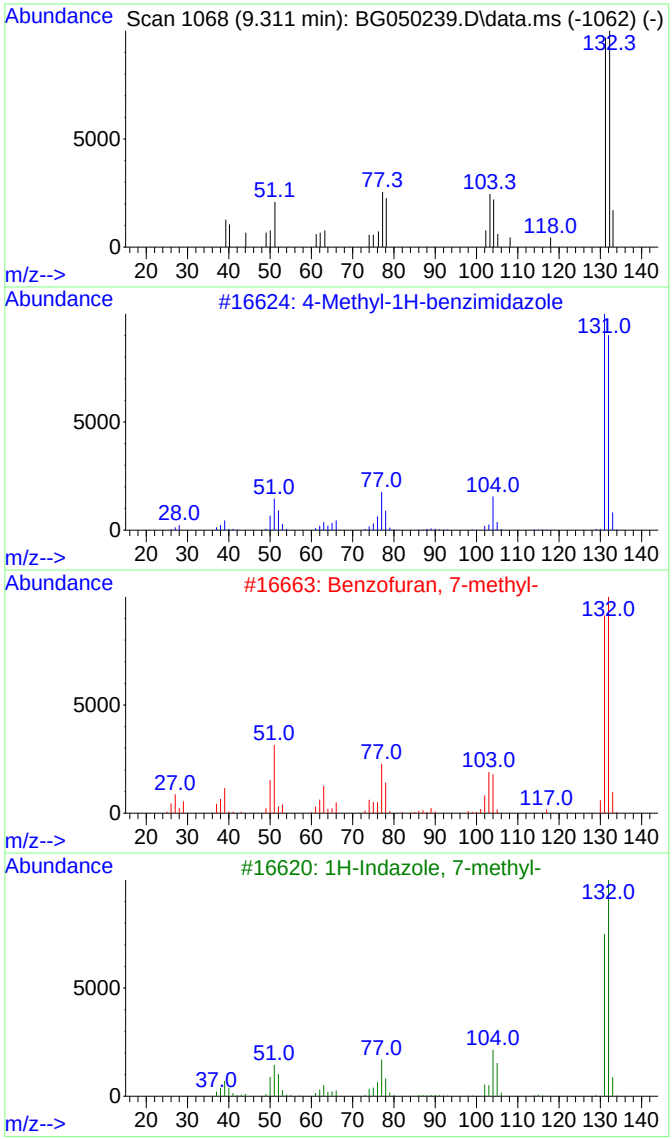
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

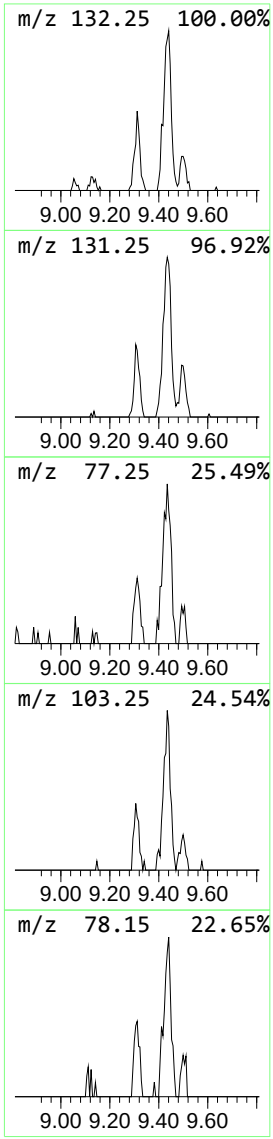
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 4-Methyl-1H-benzimidazole Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.311	2.79 ng/ul	25405	1,4-Dichlorobenzene-d4	7.949

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-1H-benzimidazole	132	C8H8N2	1000486-24-5	87
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	87
3	1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	87
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	83
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

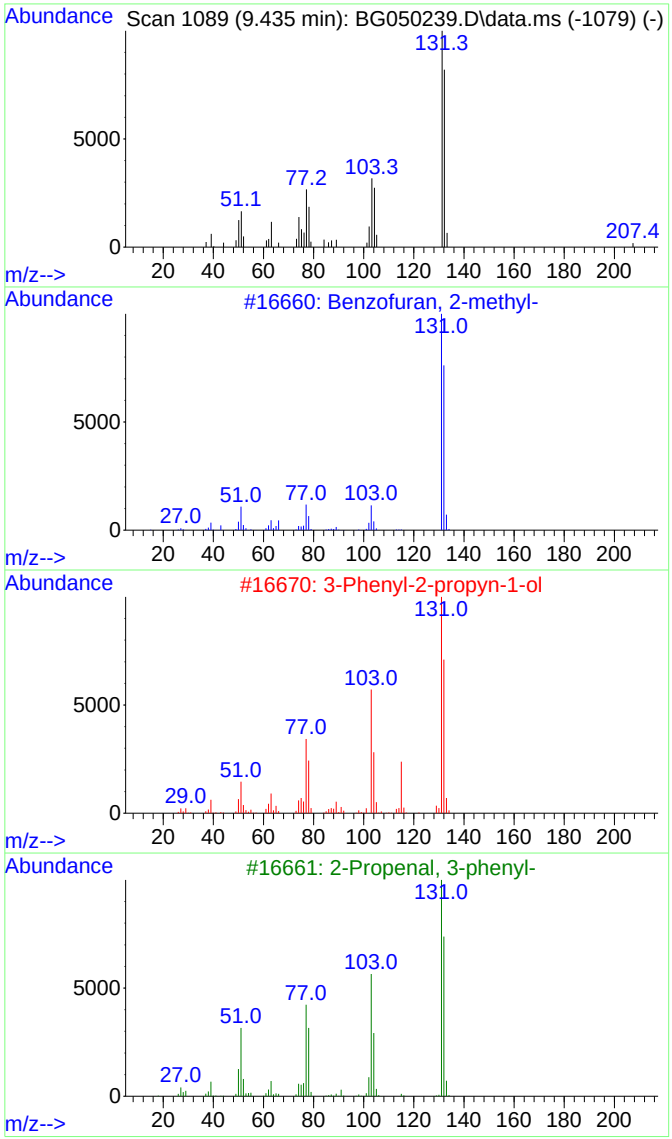
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

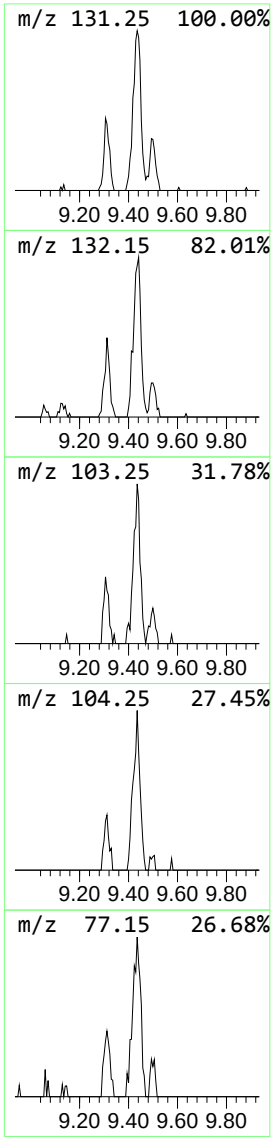
****

Peak Number 10 Benzofuran, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.435	3.60 ng/ul	78894	Naphthalene-d8	10.762

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
2	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
3	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

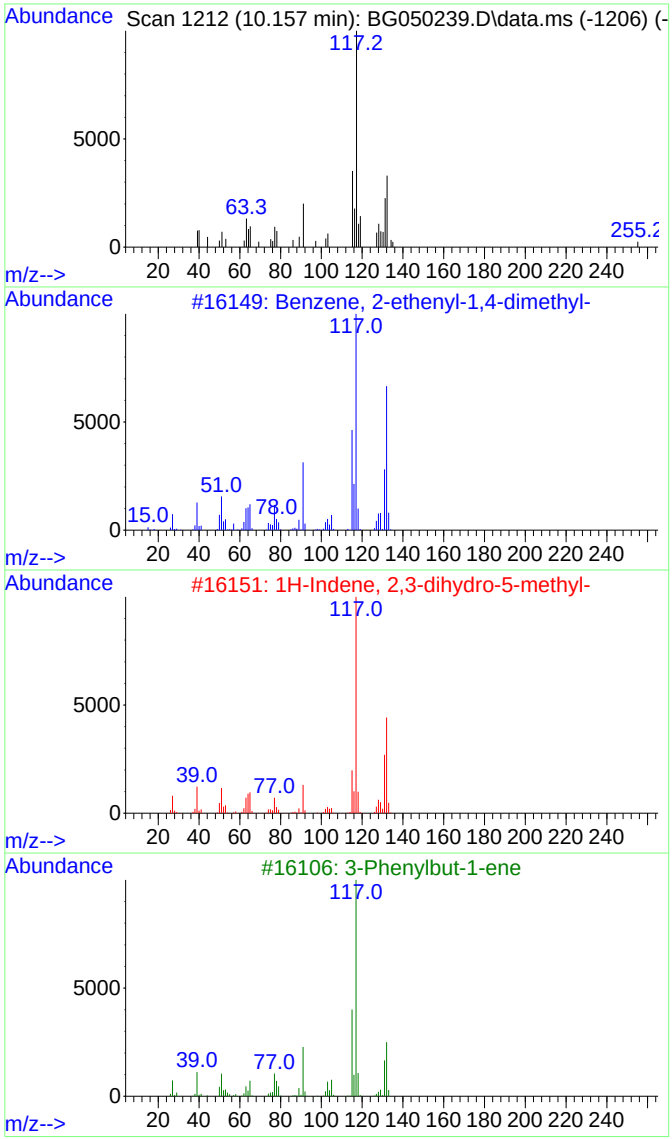
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

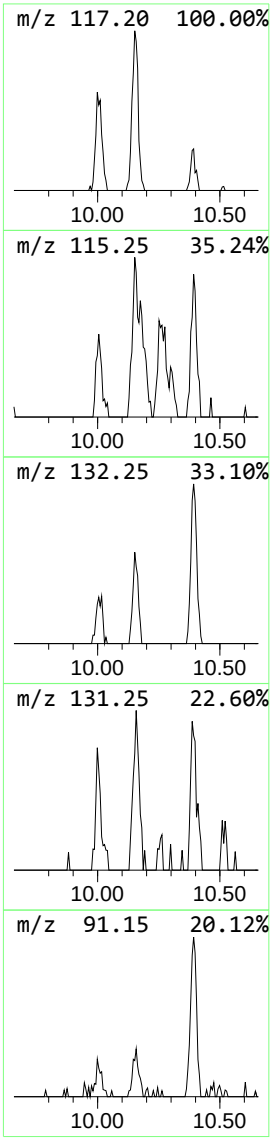
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.157	2.40 ng/ul	52665	Naphthalene-d8	10.762

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	83
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	81
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

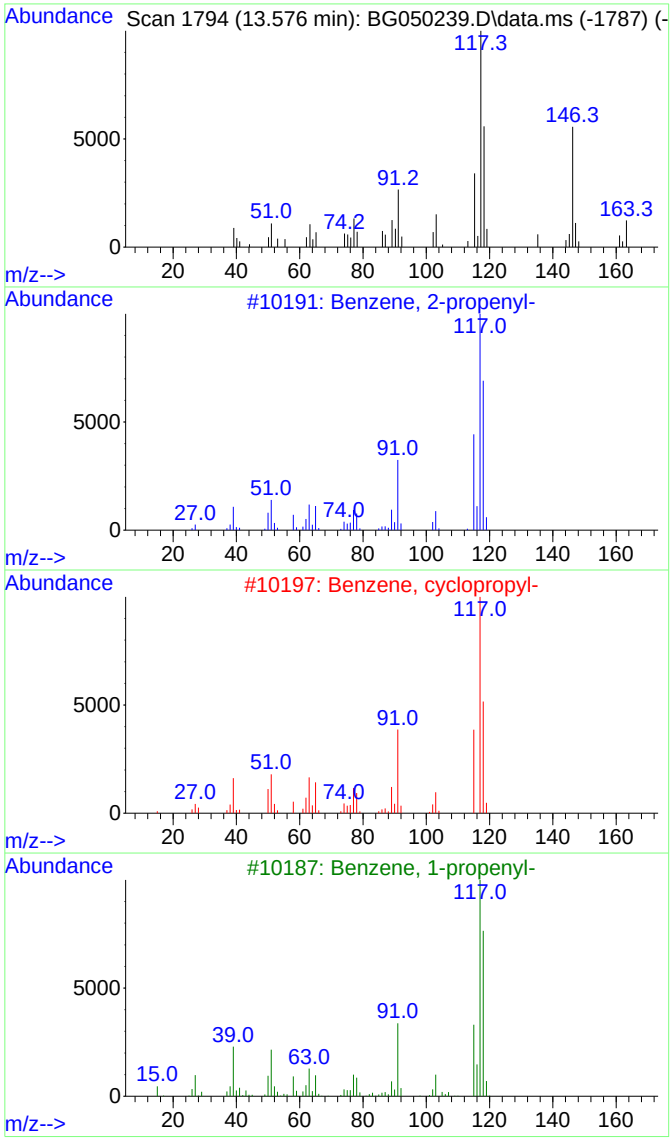
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

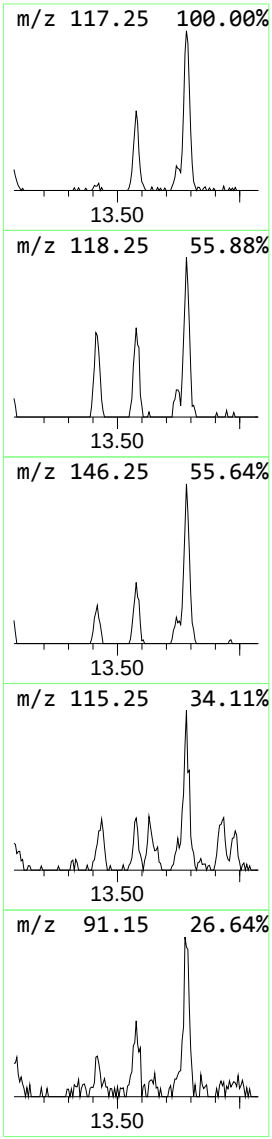
****

Peak Number 12 Benzene, 2-propenyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.576	2.16 ng/ul	47330	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	55
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBKL5DL

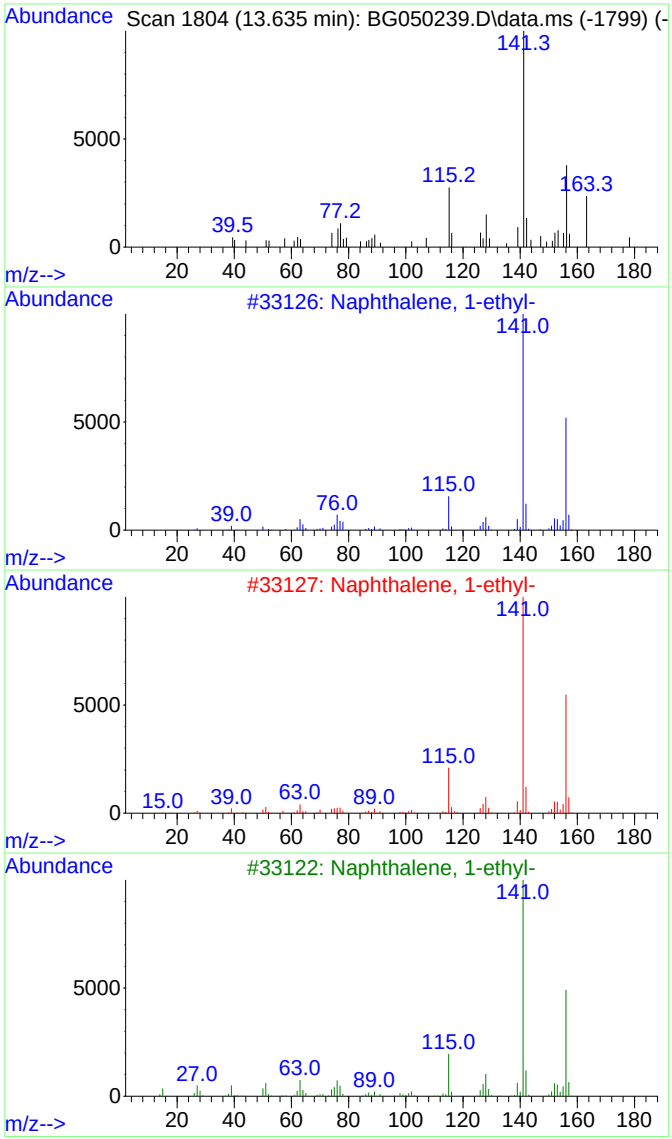
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

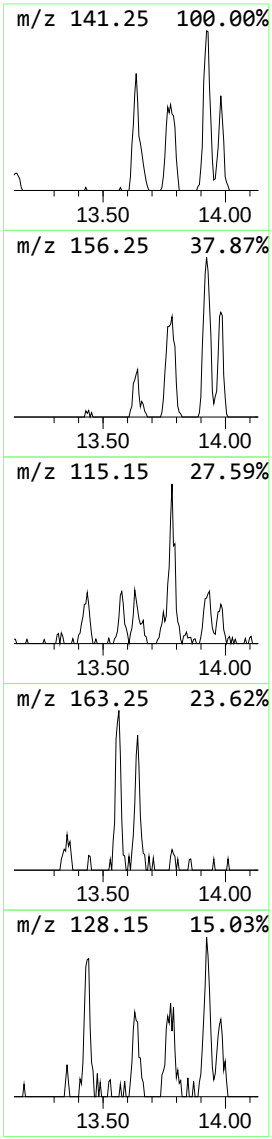
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Naphthalene, 1-ethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.635	2.39 ng/ul	52421	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	83
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	83
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	81
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	76
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

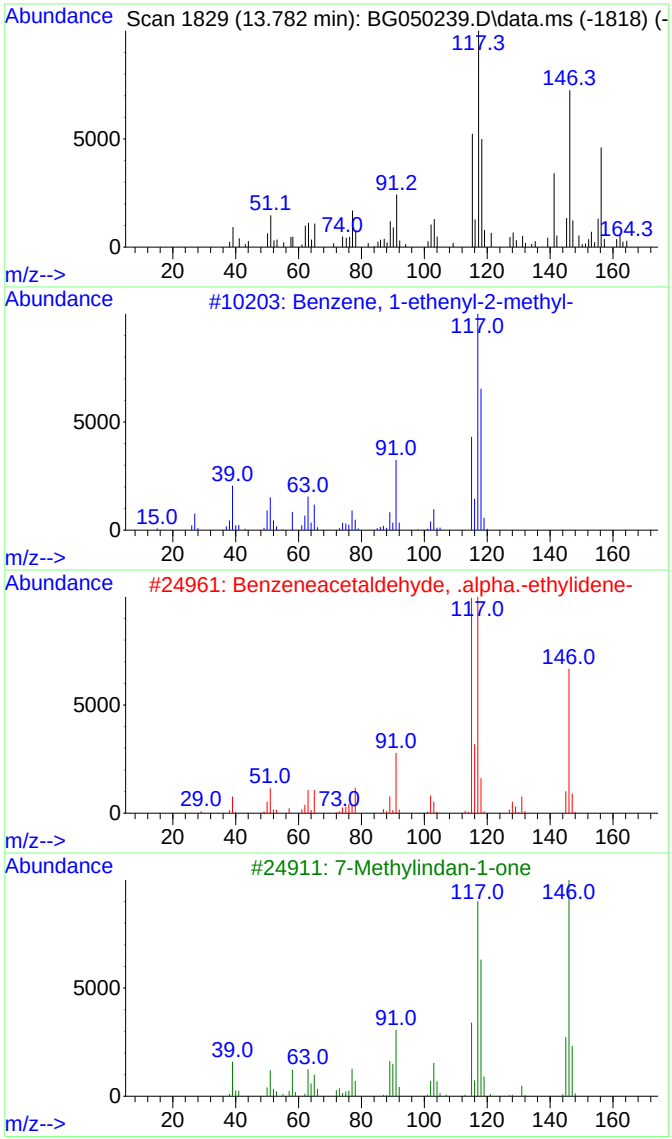
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

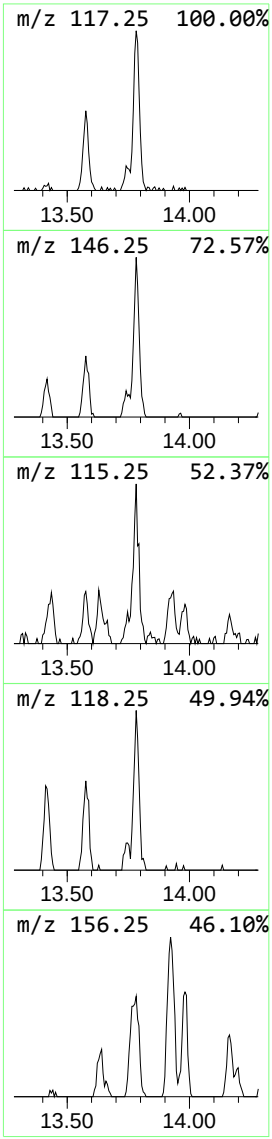
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-ethenyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.782	8.06 ng/ul	177046	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2	Benzeneacetaldehyde, .alpha.-eth...	146	C10H10O	004411-89-6	41
3	7-Methylindan-1-one	146	C10H10O	039627-61-7	38
4	Benzene, 1-pentenyl-	146	C11H14	000826-18-6	38
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

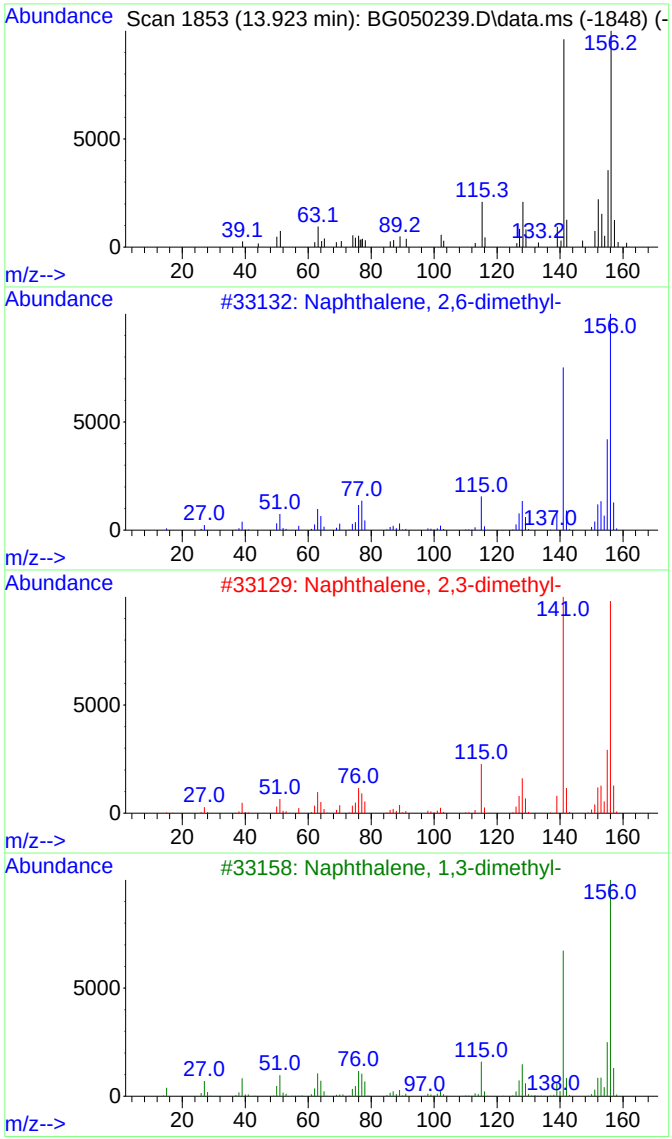
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

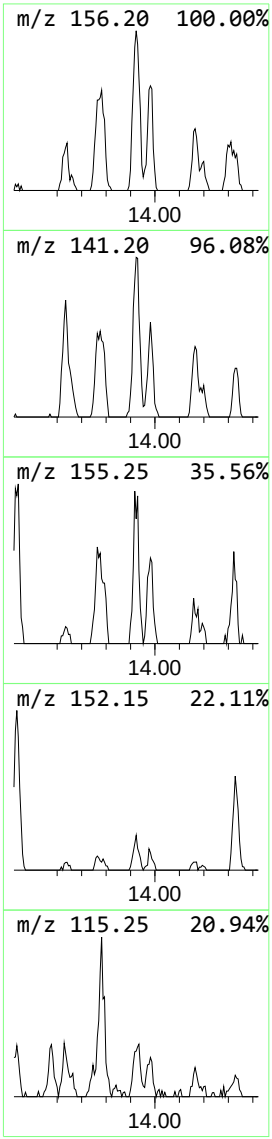
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Naphthalene, 2,6-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.923	3.46 ng/ul	75940	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

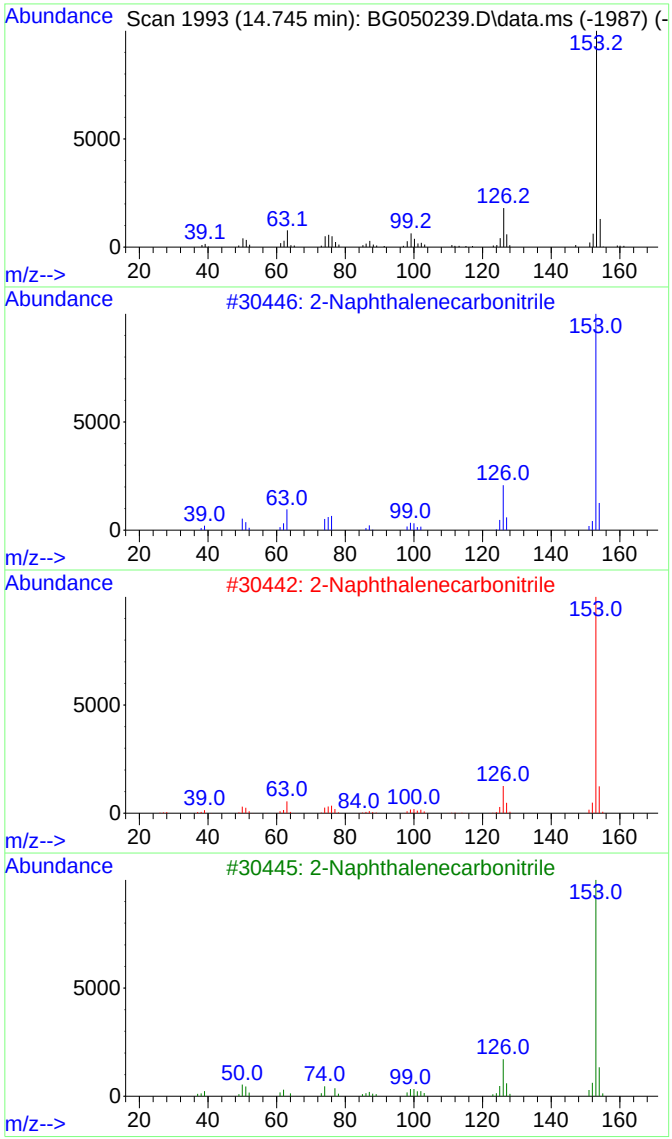
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

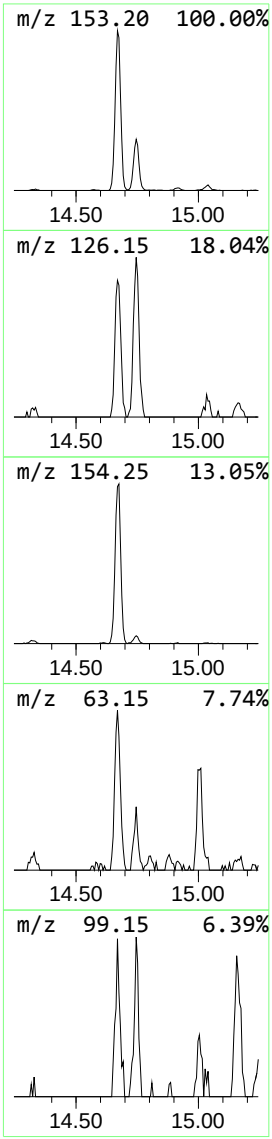
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 2-Naphthalenecarbonitrile Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.745	5.00 ng/ul	109840	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	95
2	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	94
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
5	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

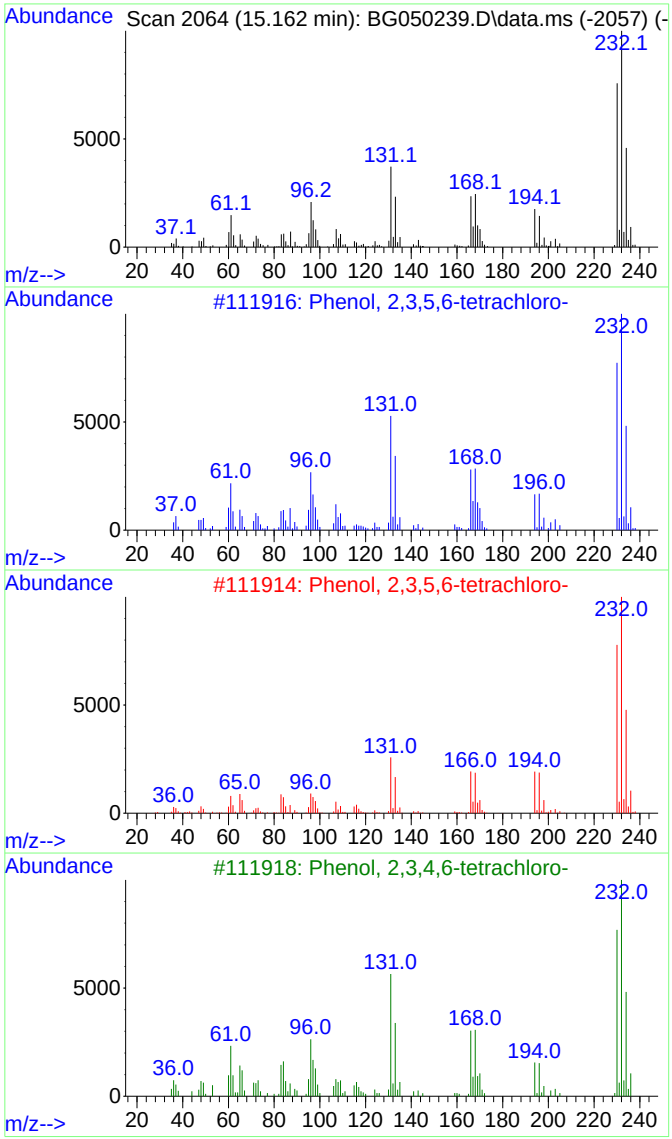
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

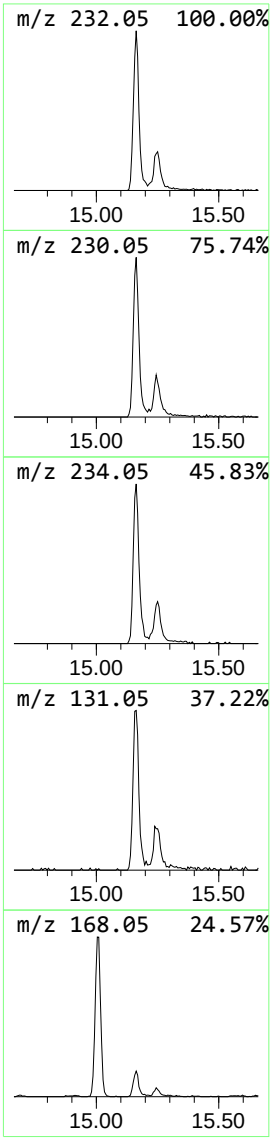
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.162	24.14 ng/ul	530180	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
2	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	96
3	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	96
4	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	95
5	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

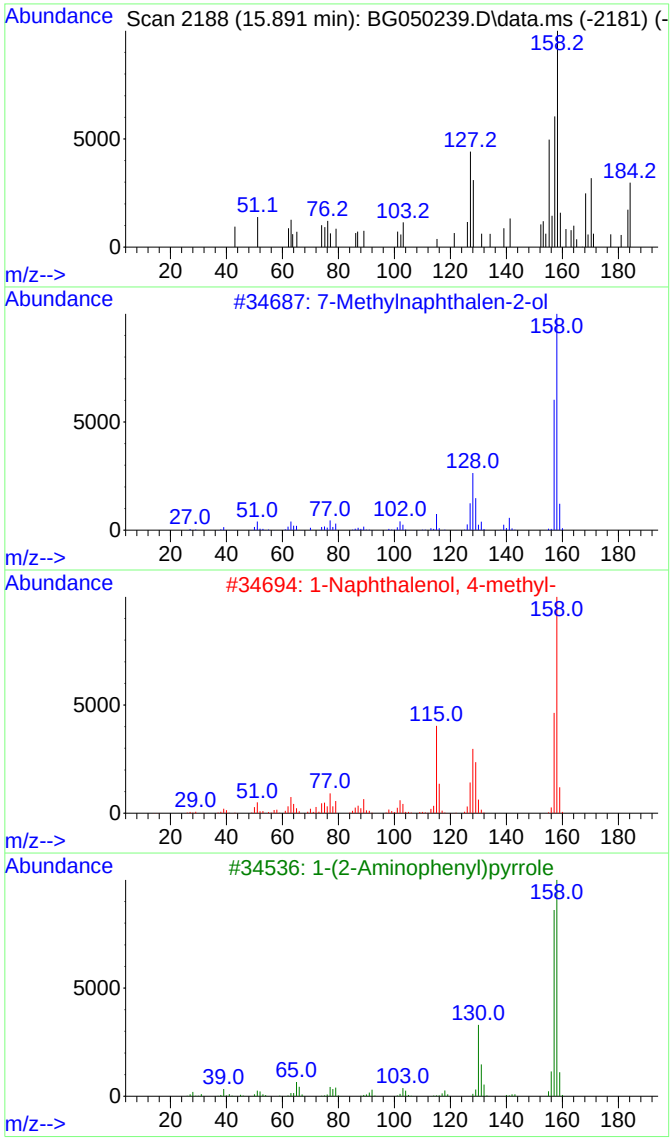
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

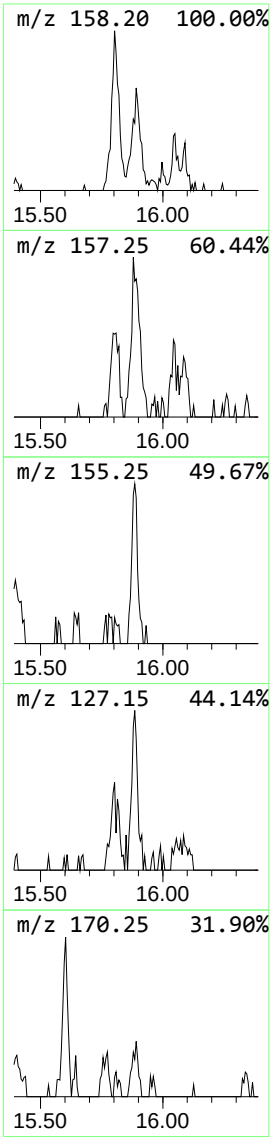
****

Peak Number 18 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.891	2.61 ng/ul	57301	Acenaphthene-d10	14.604

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methylnaphthalen-2-ol	158	C11H10O	026593-50-0	43
2	1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	43
3	1-(2-Aminophenyl)pyrrole	158	C10H10N2	006025-60-1	38
4	8-Methylquinolin-4-amine	158	C10H10N2	893762-15-7	30
5	3-Methyl-4-phenylpyrazole	158	C10H10N2	013788-84-6	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

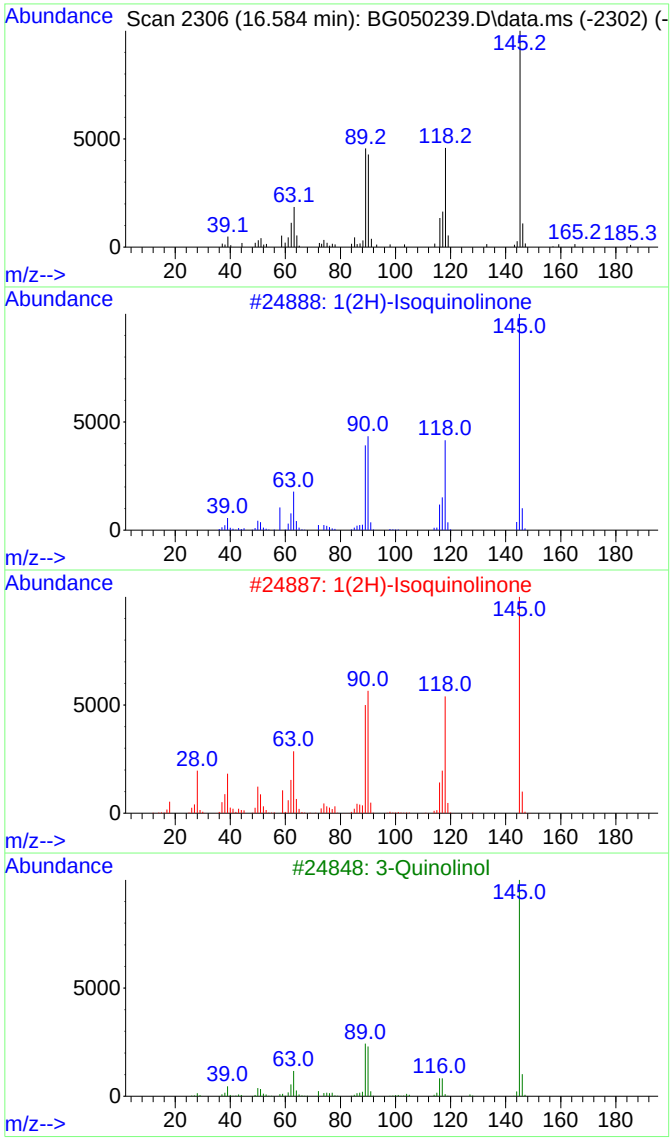
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

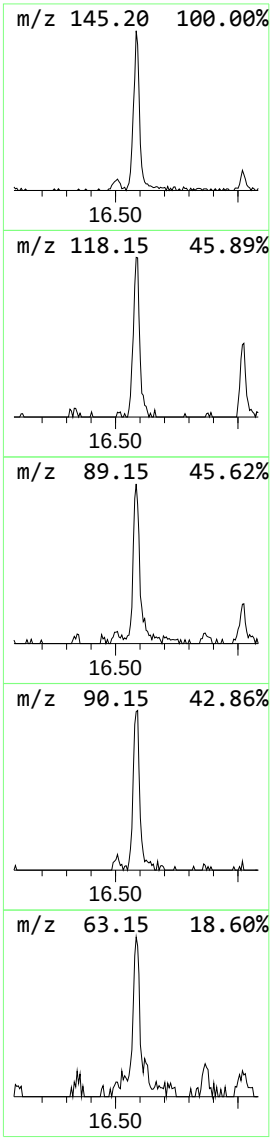
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 1(2H)-Isoquinolinone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.584	4.53 ng/ul	117800	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	95
2	1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	94
3	3-Quinolinol	145	C9H7NO	000580-18-7	81
4	1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	74
5	2-(1H-Pyrazol-3-yl)pyridine	145	C8H7N3	075415-03-1	70





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

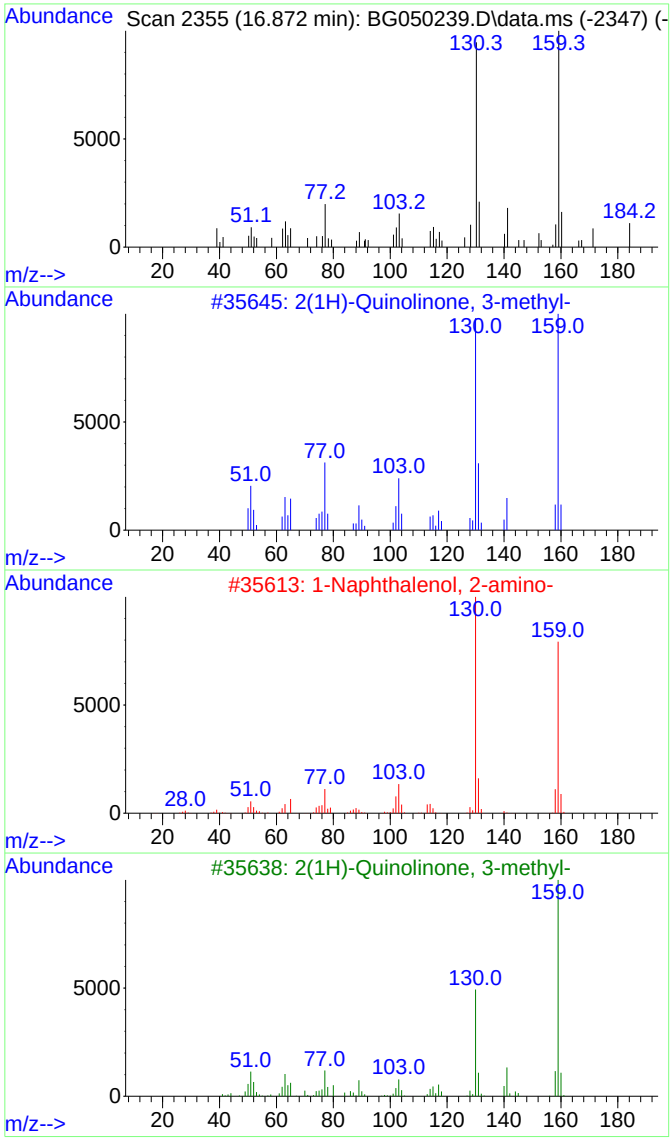
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

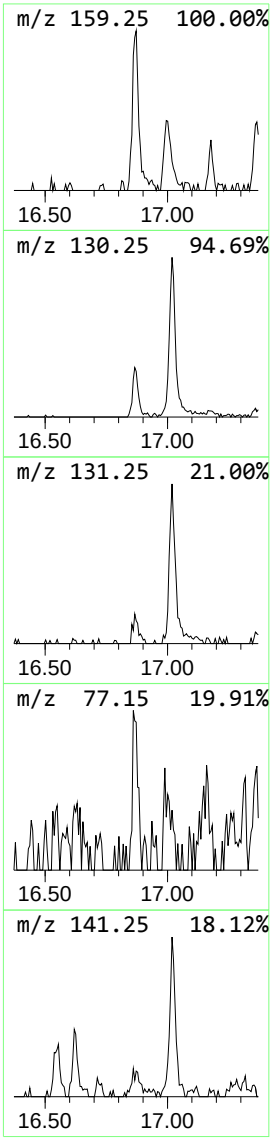
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 2(1H)-Quinolinone, 3-methyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.872	2.04 ng/ul	53087	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	90
2	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	87
3	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	87
4	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	87
5	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

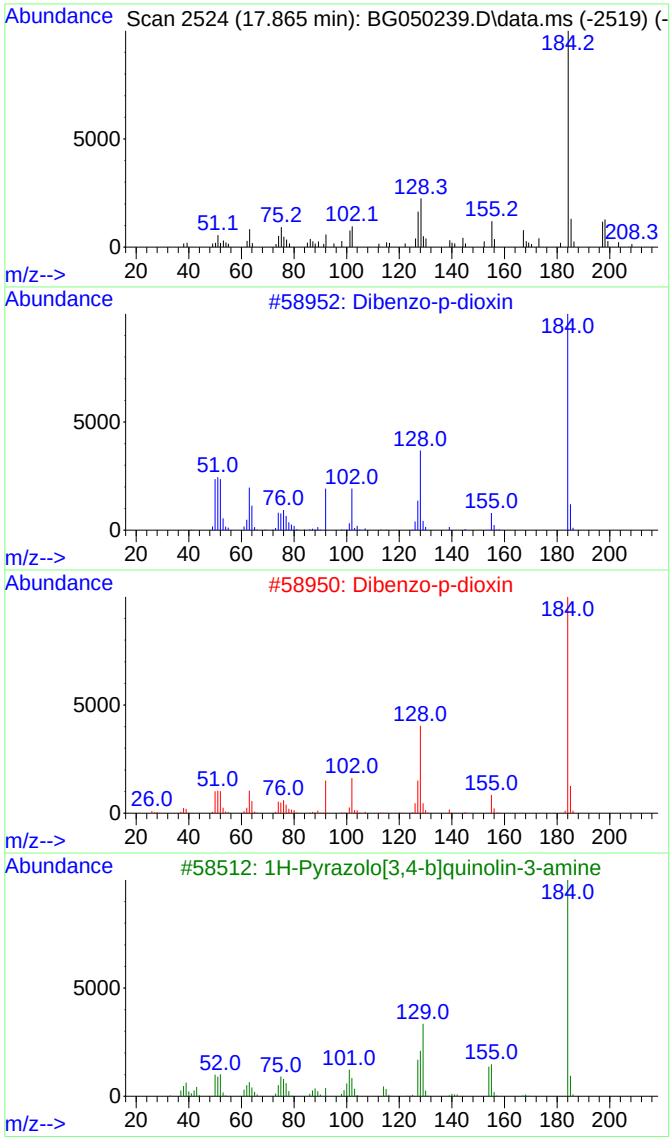
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

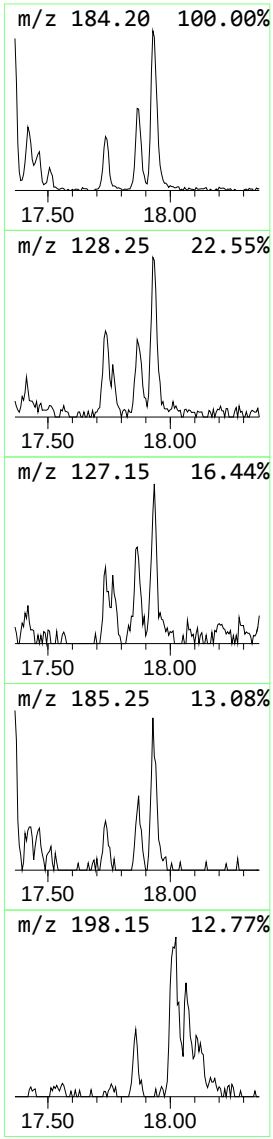
****

Peak Number 21 Dibenzo-p-dioxin Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.865	3.10 ng/ul	80472	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	83
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	83
3	1H-Pyrazolo[3,4-b]quinolin-3-amine	184	C10H8N4	1000319-54-1	74
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	74
5	2-Dibenzofuranol	184	C12H8O2	000086-77-1	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

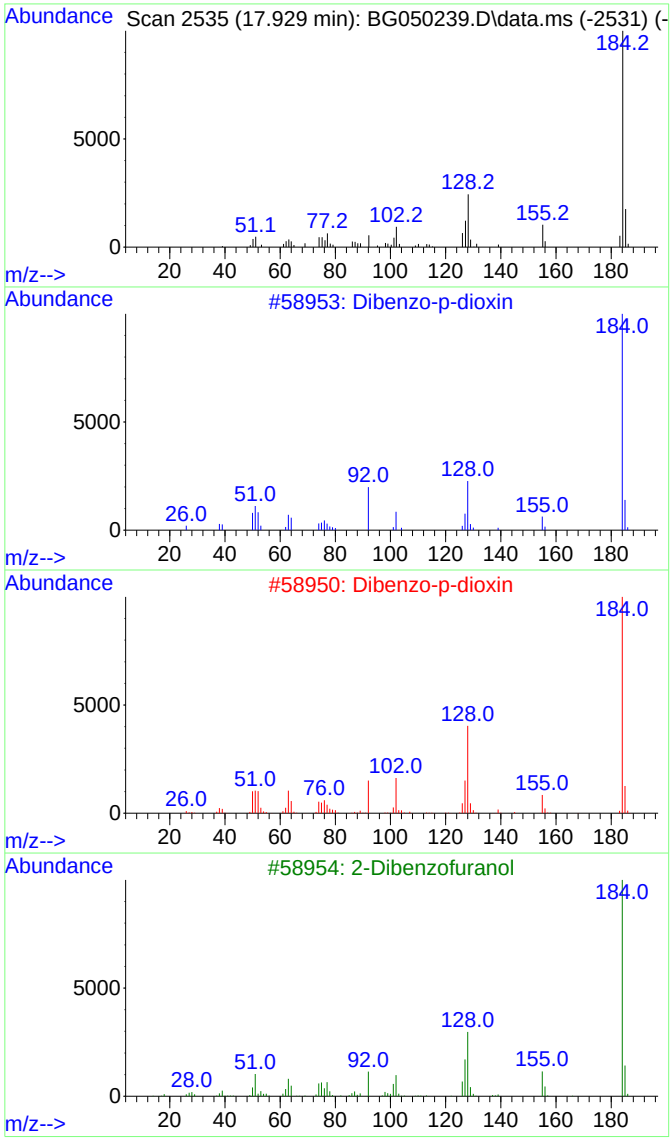
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

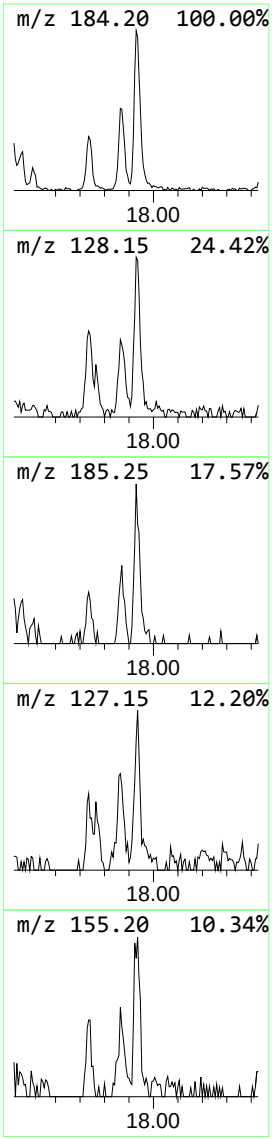
****

Peak Number 22 2-Dibenzofuranol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.929	4.06 ng/ul	105393	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	91
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	91
3	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
4	2-Dibenzofuranol	184	C12H8O2	000086-77-1	91
5	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

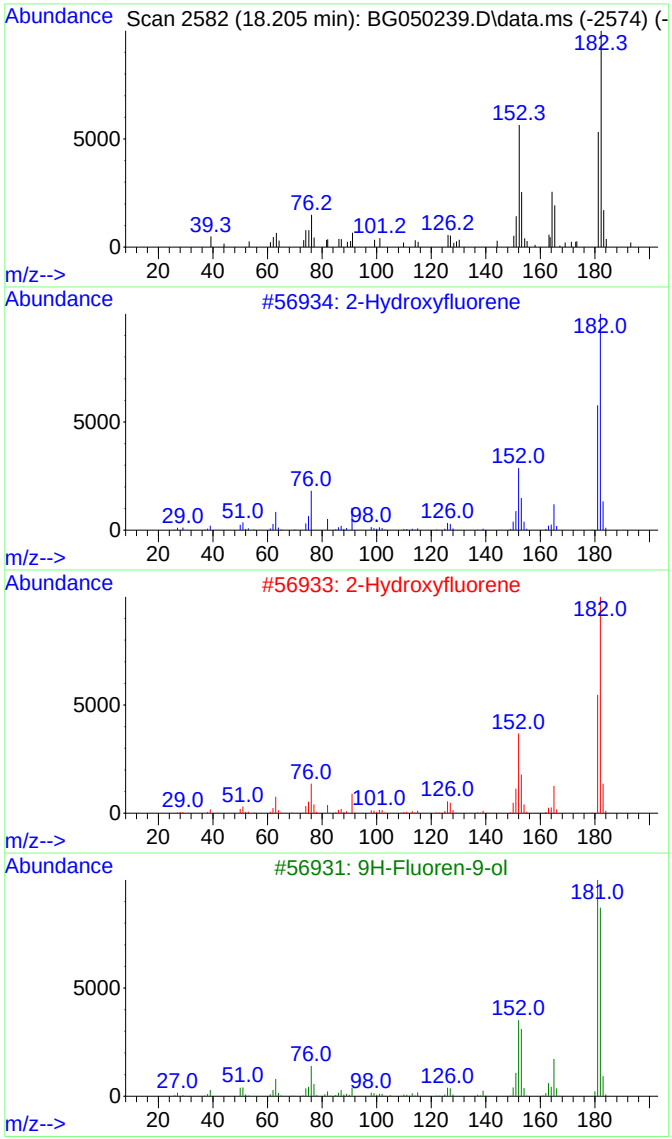
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

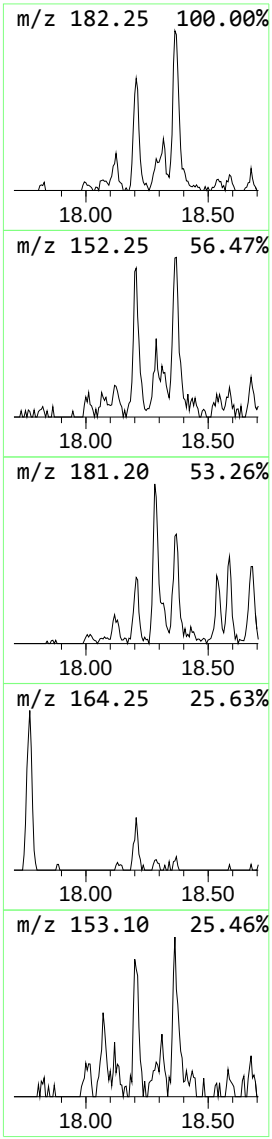
****

Peak Number 23 2-Hydroxyfluorene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.205	2.26 ng/ul	58656	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	87
2	2-Hydroxyfluorene	182	C13H10O	002443-58-5	76
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64
4	Atractylodin	182	C13H10O	055290-63-6	64
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

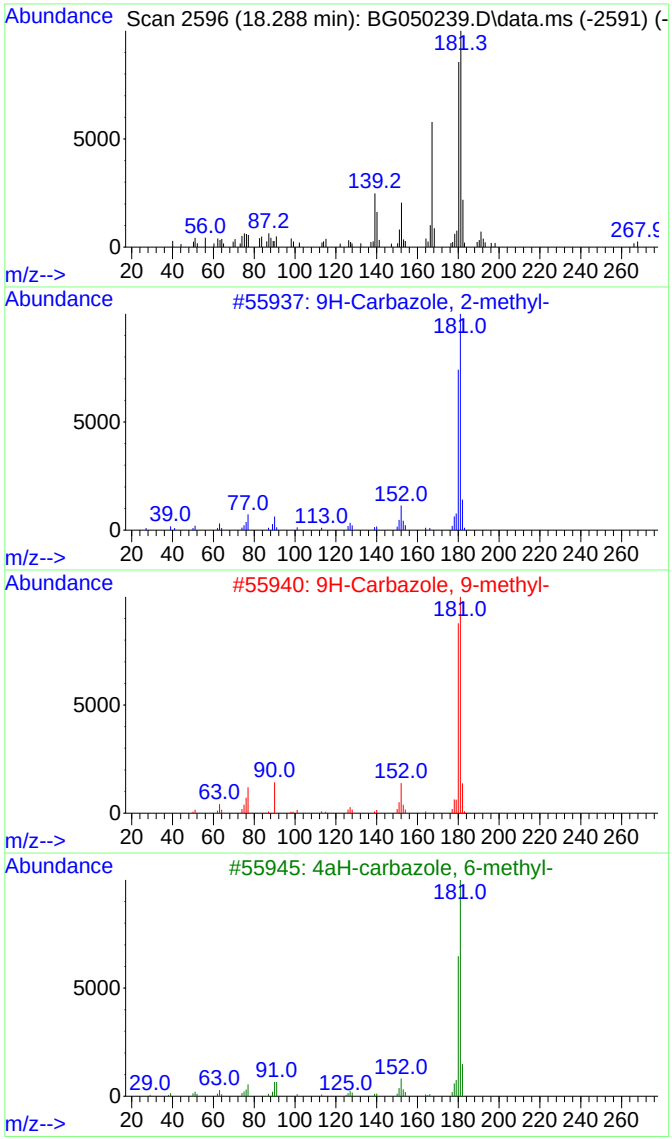
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

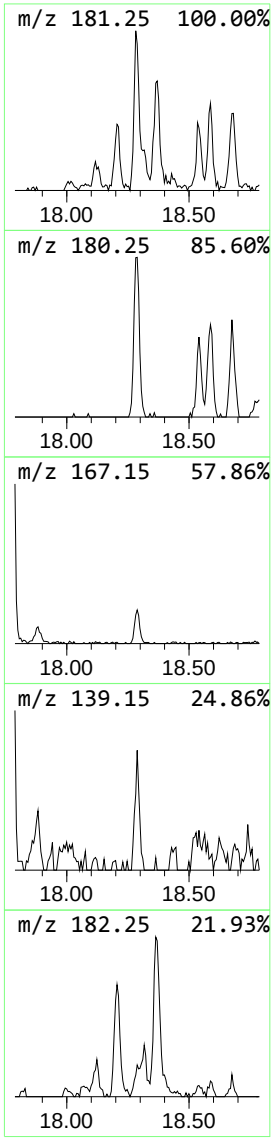
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 9H-Carbazole, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.288	3.32 ng/ul	86243	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	64
2	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	60
3	4aH-carbazole, 6-methyl-	181	C13H11N	1000404-86-1	53
4	2-Fluorenamine	181	C13H11N	000153-78-6	52
5	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

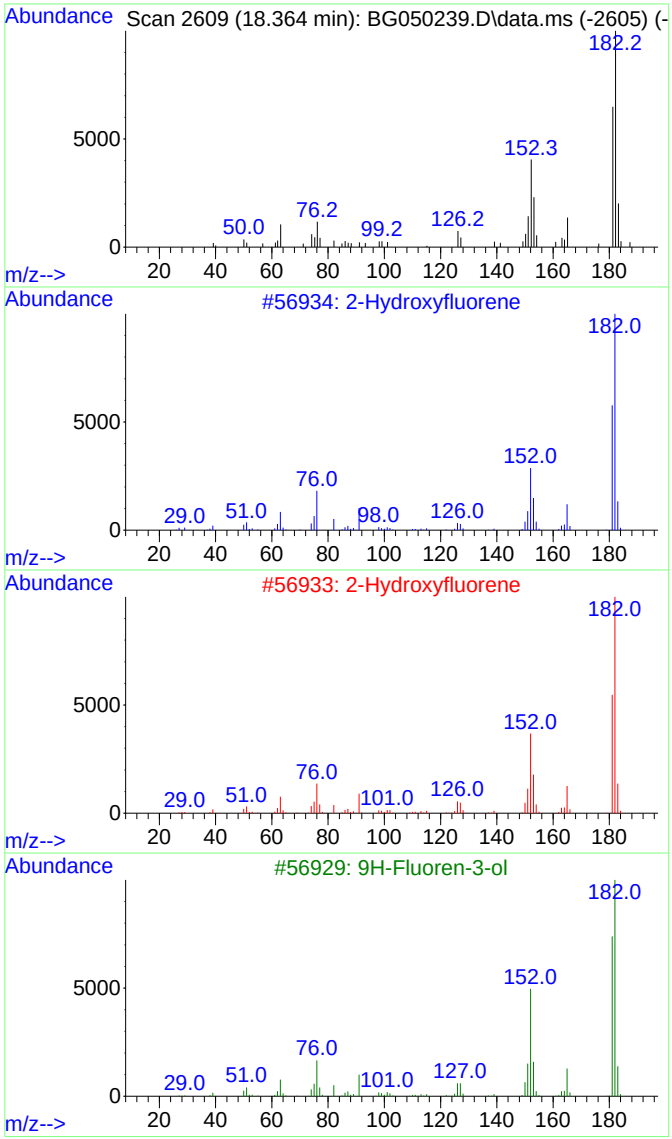
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

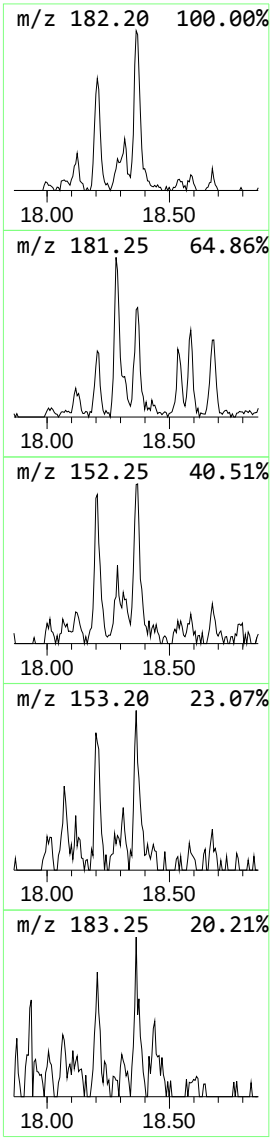
****

Peak Number 25 9H-Fluoren-3-ol Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.364	2.21 ng/ul	57374	Phenanthrene-d10	17.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hydroxyfluorene	182	C13H10O	002443-58-5	95
2	2-Hydroxyfluorene	182	C13H10O	002443-58-5	94
3	9H-Fluoren-3-ol	182	C13H10O	006344-67-8	90
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	80
5	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

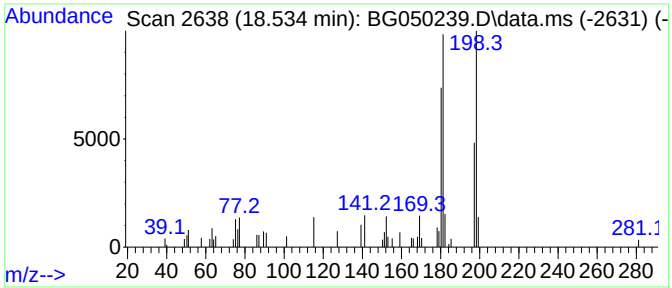
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

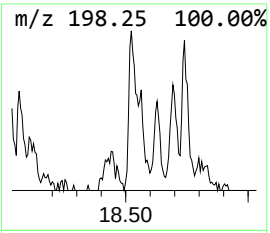
****

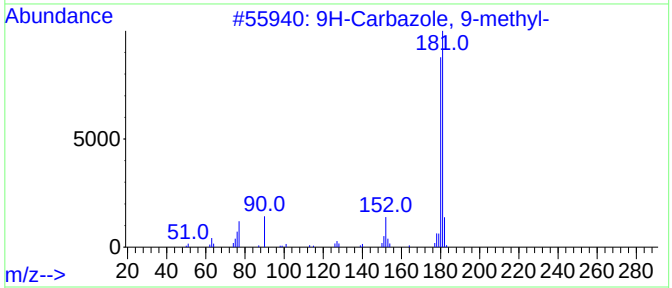
Peak Number 26 9H-Carbazole, 9-methyl- Concentration Rank 21

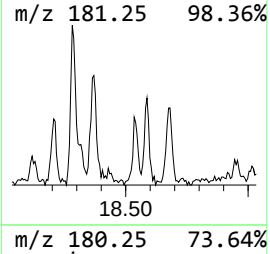
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.534	2.66 ng/ul	68998	Phenanthrene-d10	17.359

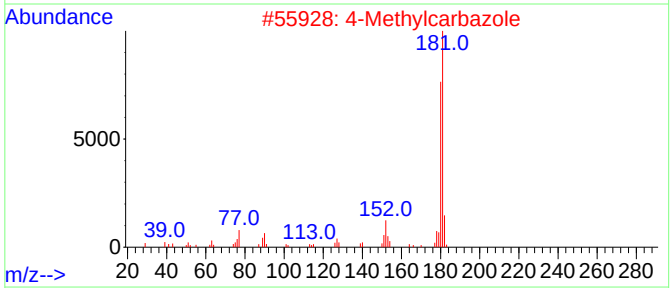
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	55
2	4-Methylcarbazole	181	C13H11N	003770-48-7	55
3	9H-Carbazole, 2-methyl-	181	C13H11N	003652-91-3	55
4	1-Methylcarbazole	181	C13H11N	006510-65-2	55
5	9H-Carbazole, 9-methyl-	181	C13H11N	001484-12-4	50

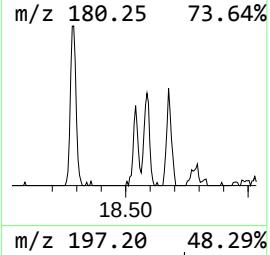


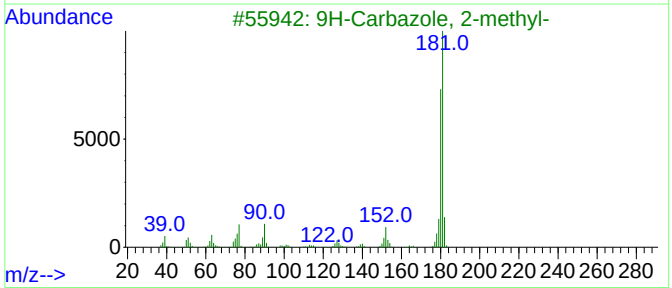


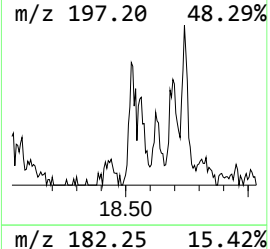


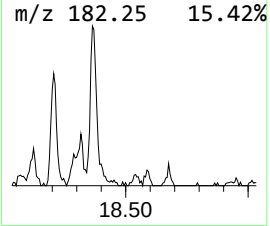












Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

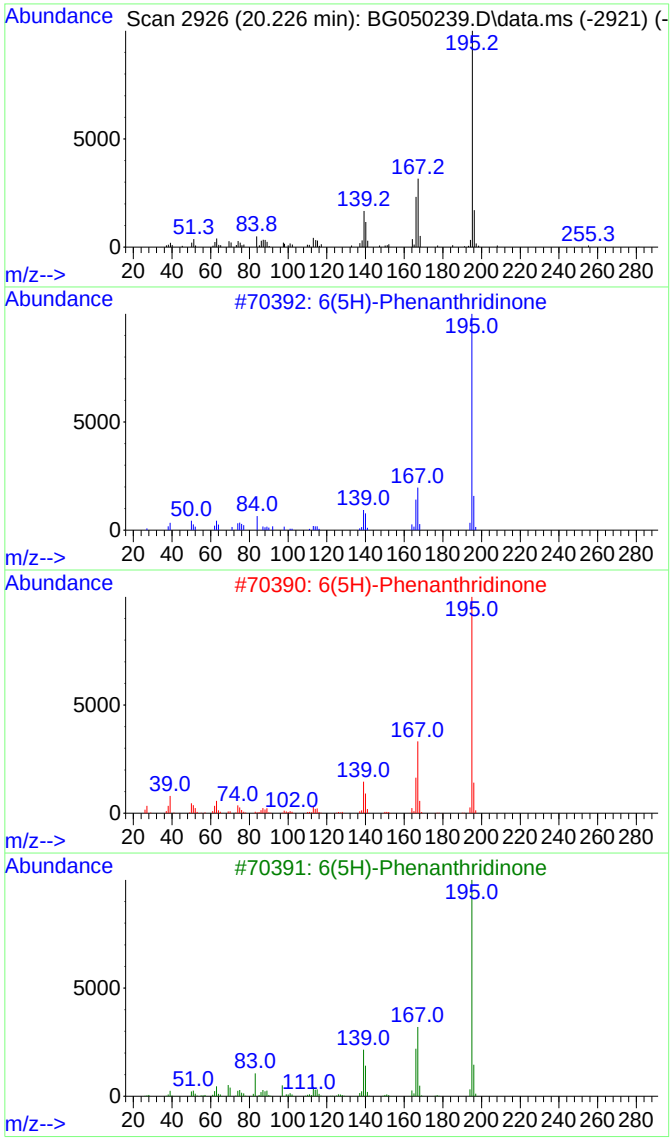
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

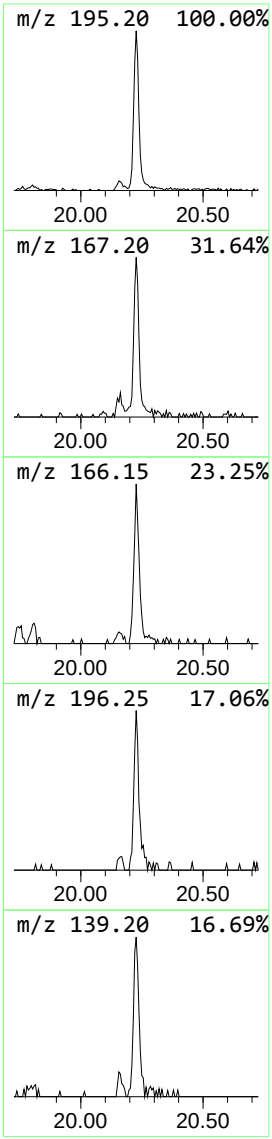
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 6(5H)-Phenanthridinone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.226	3.78 ng/ul	117912	Chrysene-d12	21.630

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	96
2	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	96
3	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	95
4	3-Acridinol	195	C13H9NO	007132-70-9	94
5	9-Acridinol	195	C13H9NO	000643-62-9	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

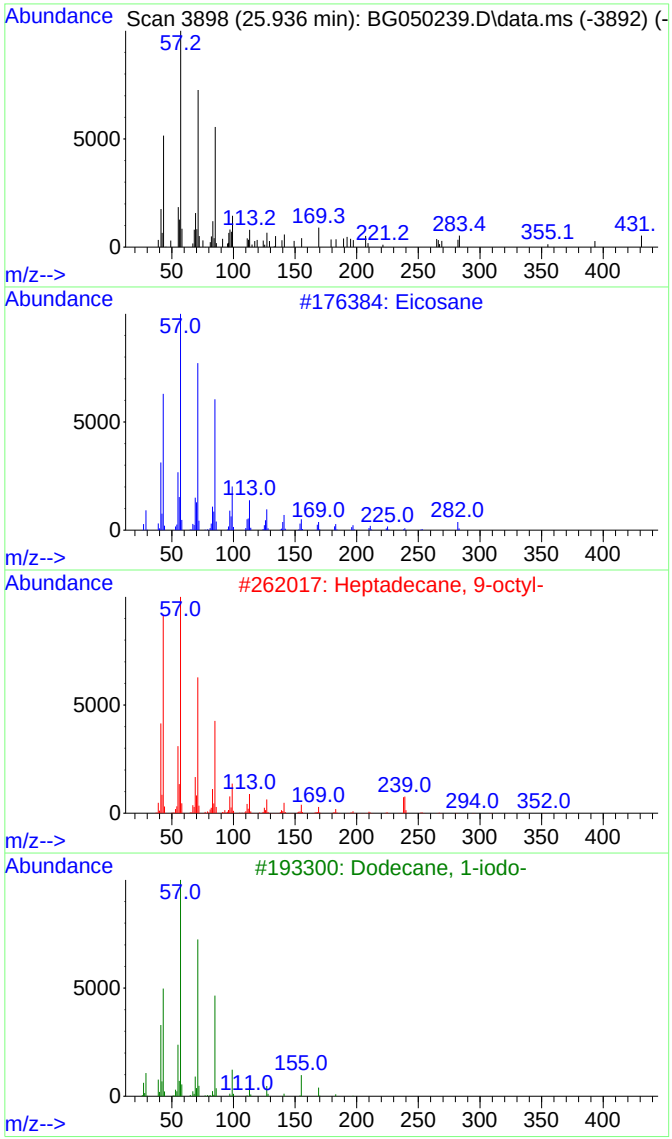
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION

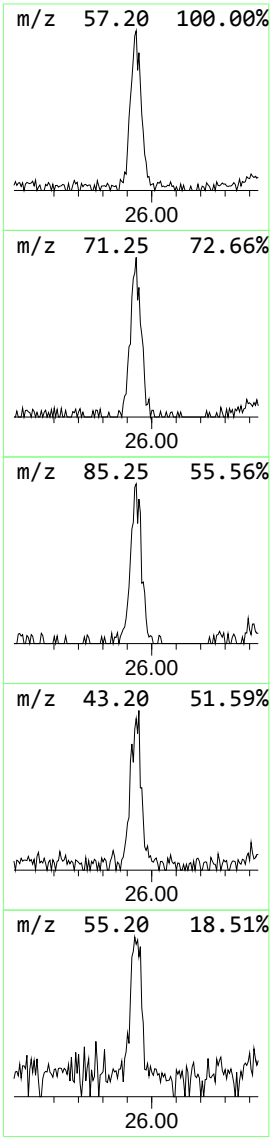
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.936	3.09 ng/ul	106840	Perylene-d12	24.843

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	81
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	74
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
4	Tetracosane	338	C24H50	000646-31-1	72
5	Pentacosane	352	C25H52	000629-99-2	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050239.D
Acq On : 9 Sep 2021 21:28
Operator : CG/JU
Sample : M3608-19DL 10X
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBKL5DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.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
Benzene, 1,3-di...	5.552	9.2	ng/ul	83953	1	7.949	182339	20.0
Benzene, 1,2,4-...	7.584	4.1	ng/ul	37354	1	7.949	182339	20.0
Benzofuran	7.661	11.6	ng/ul	105623	1	7.949	182339	20.0
Indane	8.319	17.3	ng/ul	157896	1	7.949	182339	20.0
Benzene, 1-prop...	8.483	64.7	ng/ul	589468	1	7.949	182339	20.0
4-Methyl-1H-ben...	9.311	2.8	ng/ul	25405	1	7.949	182339	20.0
Benzofuran, 2-m...	9.435	3.6	ng/ul	78894	2	10.762	437996	20.0
Benzene, 2-ethe...	10.157	2.4	ng/ul	52665	2	10.762	437996	20.0
Benzene, 2-prop...	13.576	2.2	ng/ul	47330	3	14.604	439200	20.0
Naphthalene, 1-...	13.635	2.4	ng/ul	52421	3	14.604	439200	20.0
Benzene, 1-ethe...	13.782	8.1	ng/ul	177046	3	14.604	439200	20.0
Naphthalene, 2,...	13.923	3.5	ng/ul	75940	3	14.604	439200	20.0
2-Naphthaleneca...	14.745	5.0	ng/ul	109840	3	14.604	439200	20.0
Phenol, 2,3,5,6...	15.162	24.1	ng/ul	530180	3	14.604	439200	20.0
unknown-01	15.891	2.6	ng/ul	57301	3	14.604	439200	20.0
1(2H)-Isoquinol...	16.584	4.5	ng/ul	117800	4	17.359	519634	20.0
2(1H)-Quinolino...	16.872	2.0	ng/ul	53087	4	17.359	519634	20.0
Dibenzo-p-dioxin	17.865	3.1	ng/ul	80472	4	17.359	519634	20.0
2-Dibenzofuranol	17.929	4.1	ng/ul	105393	4	17.359	519634	20.0
2-Hydroxyfluorene	18.205	2.3	ng/ul	58656	4	17.359	519634	20.0
9H-Carbazole, 2...	18.288	3.3	ng/ul	86243	4	17.359	519634	20.0
9H-Fluoren-3-ol	18.364	2.2	ng/ul	57374	4	17.359	519634	20.0
9H-Carbazole, 9...	18.534	2.7	ng/ul	68998	4	17.359	519634	20.0
6(5H)-Phenanthr...	20.226	3.8	ng/ul	117912	5	21.630	624684	20.0
(DEL) Alkane: S...	25.936	3.1	ng/ul	106840	6	24.843	692556	20.0