LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110419\
Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

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\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.069	10	14	20	rVB	1046135	1057103	15.75%	2.189%
2	7.598	778	784	795	rVB	1811032	2855541	42.54%	5.914%
3	8.704	970	972	977	rVB5	19554	19108	0.28%	0.040%
4	8.840	992	995	1007	rVB3	30646	54048	0.81%	0.112%
5	9.792	1153	1157	1162	rVB5	11345	21538	0.32%	0.045%
6	10.375	1248	1256	1263	rBV	2173850	3731220	55.58%	7.727%
7	10.428	1263	1265	1274	rVB	62457	83105	1.24%	0.172%
8	10.510	1276	1279	1281	rBV3	7208	8712	0.13%	0.018%
9	10.792	1324	1327	1331	rVB5	7924	10238	0.15%	0.021%
10	10.857	1334	1338	1345	rVB6	6579	11427	0.17%	0.024%
11	11.075	1368	1375	1379	rBV8	10373	22922	0.34%	0.047%
12	11.145	1384	1387	1391	rBV4	6583	7588	0.11%	0.016%
13	11.434	1431	1436	1438	rBV4	10507	14920	0.22%	0.031%
14	11.492	1444	1446	1450	rVB3	6005	7287	0.11%	0.015%
15	11.681	1473	1478	1480	rBV4	7654	10197	0.15%	0.021%
16	11.833	1501	1504	1509	rVB6	7772	10364	0.15%	0.021%
17	12.051	1536	1541	1547	rVB	25042	37285	0.56%	0.077%
18	12.275	1574	1579	1584	rBV3	20263	32014	0.48%	0.066%
19	12.675	1643	1647	1648	rBV2	6406	7489	0.11%	0.016%
20	12.745	1655	1659	1664	rVB7	8272	10725	0.16%	0.022%
21	12.810	1664	1670	1673	rBV4	9667	18287	0.27%	0.038%
22	13.039	1703	1709	1712	rVV6	20174	36191	0.54%	0.075%
23	13.092	1712	1718	1726	rVV9	20678	48389	0.72%	0.100%
24	13.180	1730	1733	1736	rBV5	6826	8528	0.13%	0.018%
25	13.280	1742	1750	1752	rBV8	16817	25191	0.38%	0.052%
26	13.422	1771	1774	1780	rVB6	10999	20069	0.30%	0.042%
27	13.575	1796	1800	1806	rBV8	19173	29550	0.44%	0.061%
28	13.622	1806	1808	1811	rVV4	10367	13435	0.20%	0.028%
29	13.675	1814	1817	1822	rVB6	12594	18333	0.27%	0.038%
30	13.763	1827	1832	1836	rVV6	14664	23158	0.34%	0.048%
31	13.804	1836	1839	1850	rVB8	17663	37110	0.55%	0.077%
32	13.904	1853	1856	1859	rBV4	7815	9272	0.14%	0.019%
33	13.969	1864	1867	1872	rBV5	16079	30809	0.46%	0.064%
34	14.051	1877	1881	1883	rVV5	9209	13084	0.19%	0.027%

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Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

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\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

35	14.086	1884	1887	1891	rVB5	12292	12920	0.19%	0.027%
36	14.175	1896	1902	1906	rBV8	22172	43038	0.64%	0.089%
37	14.251	1906	1915	1928	rBV	2999008	4400754	65.56%	9.114%
38	14.345	1928	1931	1936	rVV7	19418	32358	0.48%	0.067%
39	14.386	1936	1938	1940	rVB3	7642	7108	0.11%	0.015%
40	14.463	1948	1951	1953	rBV4	8547	10094	0.15%	0.021%
41	14.574	1967	1970	1975	rVV6	15260	20929	0.31%	0.043%
42	14.651	1979	1983	1991	rVV2	34387	66731	0.99%	0.138%
43	14.733	1995	1997	2000	rVB4	11688	9407	0.14%	0.019%
44	14.804	2006	2009	2013	rVB5	11973	17810	0.27%	0.037%
45	14.880	2017	2022	2023	rBV5	13357	13317	0.20%	0.028%
46	14.922	2026	2029	2034	rVV7	10012	15412	0.23%	0.032%
47	15.051	2045	2051	2055	rBV8	20566	35702	0.53%	0.074%
48	15.139	2062	2066	2070	rBV6	23680	41744	0.62%	0.086%
49	15.304	2089	2094	2099	rBV	60743	95207	1.42%	0.197%
50	15.427	2113	2115	2118	rBV4	7884	9837	0.15%	0.020%
51	15.469	2119	2122	2127	rVV6	17099	27668	0.41%	0.057%
52	15.545	2127	2135	2138	rVV7	26180	60711	0.90%	0.126%
53	15.757	2166	2171	2172	rBV5	23323	31592	0.47%	0.065%
54	15.998	2206	2212	2213	rBV5	42165	57737	0.86%	0.120%
55	16.316	2262	2266	2269	rBV5	41089	52171	0.78%	0.108%
56	16.539	2300	2304	2307	rBV3	56988	75375	1.12%	0.156%
57	16.886	2358	2363	2366	rBV	353959	466041	6.94%	0.965%
58	16.921	2366	2369	2373	rVB	137656	176967	2.64%	0.366%
59	16.998	2376	2382	2386	rBV	3175503	4571733	68.10%	9.468%
60	17.039	2386	2389	2396	rVB	1138382	1534221	22.85%	3.177%
61	17.180	2409	2413	2422	rVB3	224499	353289	5.26%	0.732%
62	17.457	2457	2460	2463	rBV3	65422	75601	1.13%	0.157%
63	17.604	2479	2485	2493	rBV	390863	795458	11.85%	1.647%
64	17.721	2501	2505	2513	rVB4	213409	338999	5.05%	0.702%
65	17.804	2516	2519	2523	rVB6	53493	75058	1.12%	0.155%
66	17.915	2534	2538	2543	rVB3	201918	340902	5.08%	0.706%
67	18.074	2560	2565	2572	rBV3	1004415	1473610	21.95%	3.052%
68	18.139	2572	2576	2579	rVV	596796	778920	11.60%	1.613%
69	18.180	2579	2583	2588	rVB	429682	552781	8.23%	1.145%
70	18.362	2610	2614	2619	rBV2	716699	1058632	15.77%	2.192%
71	18.410	2619	2622	2626	rVB	223793	265141	3.95%	0.549%

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ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

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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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

72	18.504	2635	2638	2641	rBV2	226275	259845	3.87%	0.538%
73	18.539	2641	2644	2651	rVB	401886	508011	7.57%	1.052%
74	18.639	2658	2661	2664	rVB3	151591	165362	2.46%	0.342%
75	18.857	2694	2698	2701	rBV3	248024	335817	5.00%	0.695%
76	18.904	2702	2706	2709	rVV	791504	1107223	16.49%	2.293%
77	18.939	2709	2712	2718	rVB2	835921	1187692	17.69%	2.460%
78	19.062	2728	2733	2740	rVB	2083845	2795495	41.64%	5.789%
79	19.215	2755	2759	2765	rBV2	387908	587423	8.75%	1.217%
80	19.280	2768	2770	2773	rVB	173510	170223	2.54%	0.353%
81	19.398	2787	2790	2793	rBV	188876	282573	4.21%	0.585%
82	19.662	2832	2835	2840	rVB	385773	466250	6.95%	0.966%
83	19.980	2886	2889	2892	rBV	340739	375988	5.60%	0.779%
84	21.192	3090	3095	3099	rBV	4235038	5984540	89.15%	12.394%
85	21.227	3099	3101	3108	rVB	834325	1007577	15.01%	2.087%
86	23.380	3462	3467	3479	rVB	3569936	6712915	100.00%	13.902%

Sum of corrected areas: 48286146

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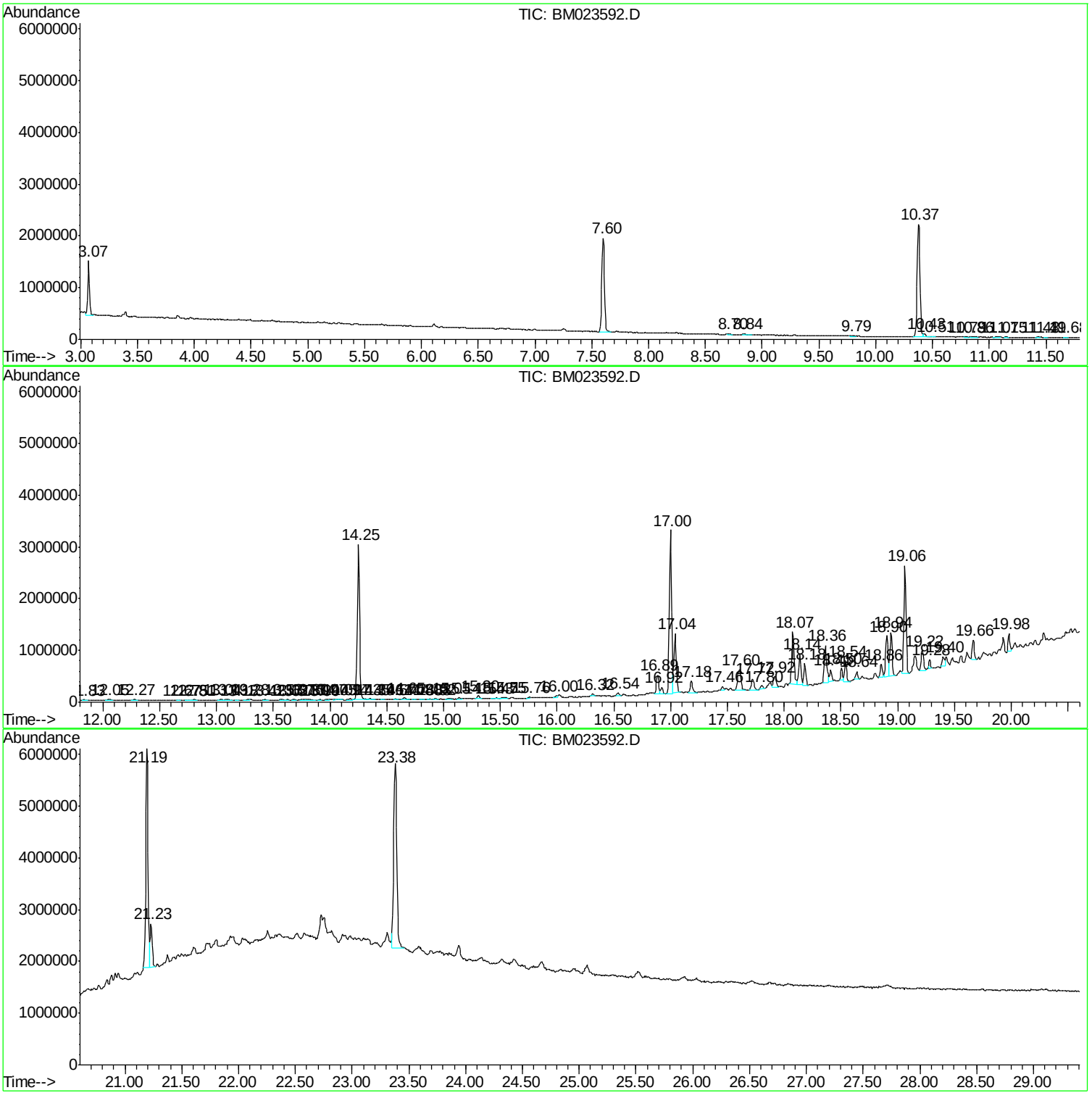
Instrument :
BNA_M
ClientSampleId :
BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

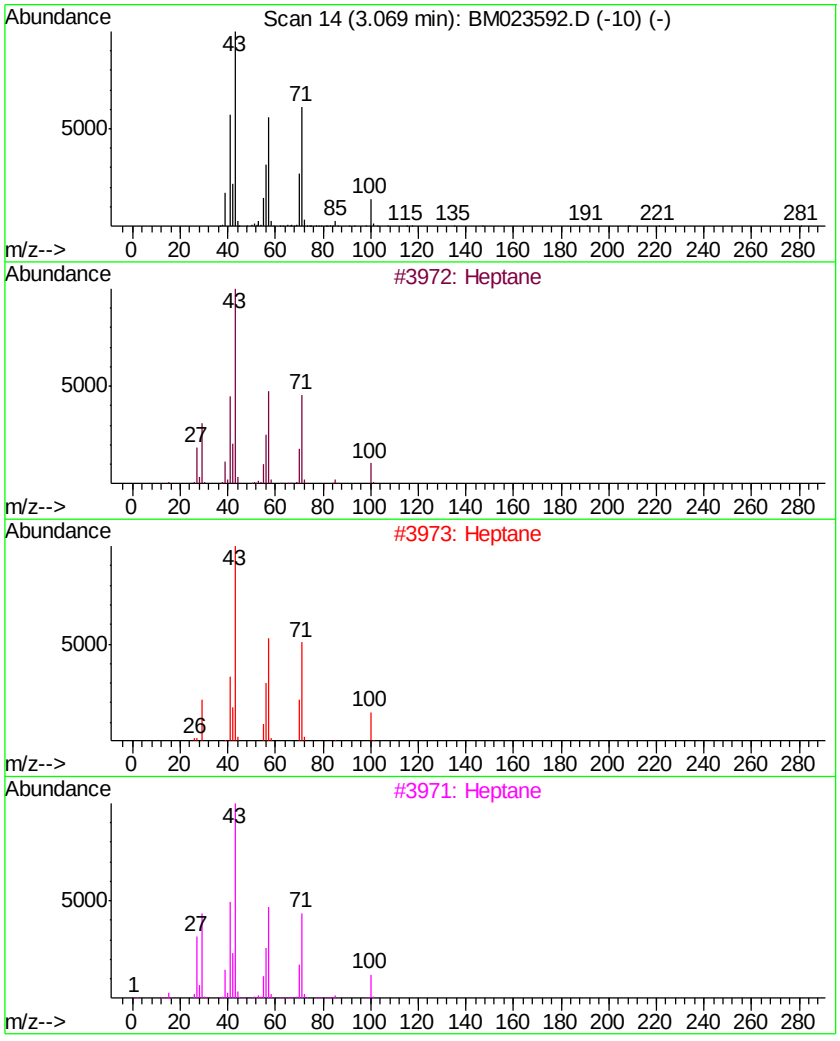
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

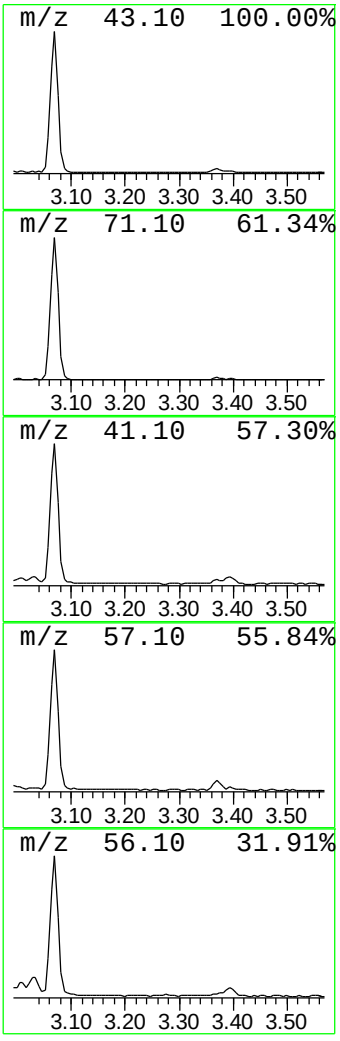
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.07	7.40 ng/ul	1057100	1,4-Dichlorobenzene-d4	7.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	94
2	Heptane	100	C7H16	000142-82-5	91
3	Heptane	100	C7H16	000142-82-5	90
4	Heptane	100	C7H16	000142-82-5	64
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	40





Library Search Compound Report

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Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

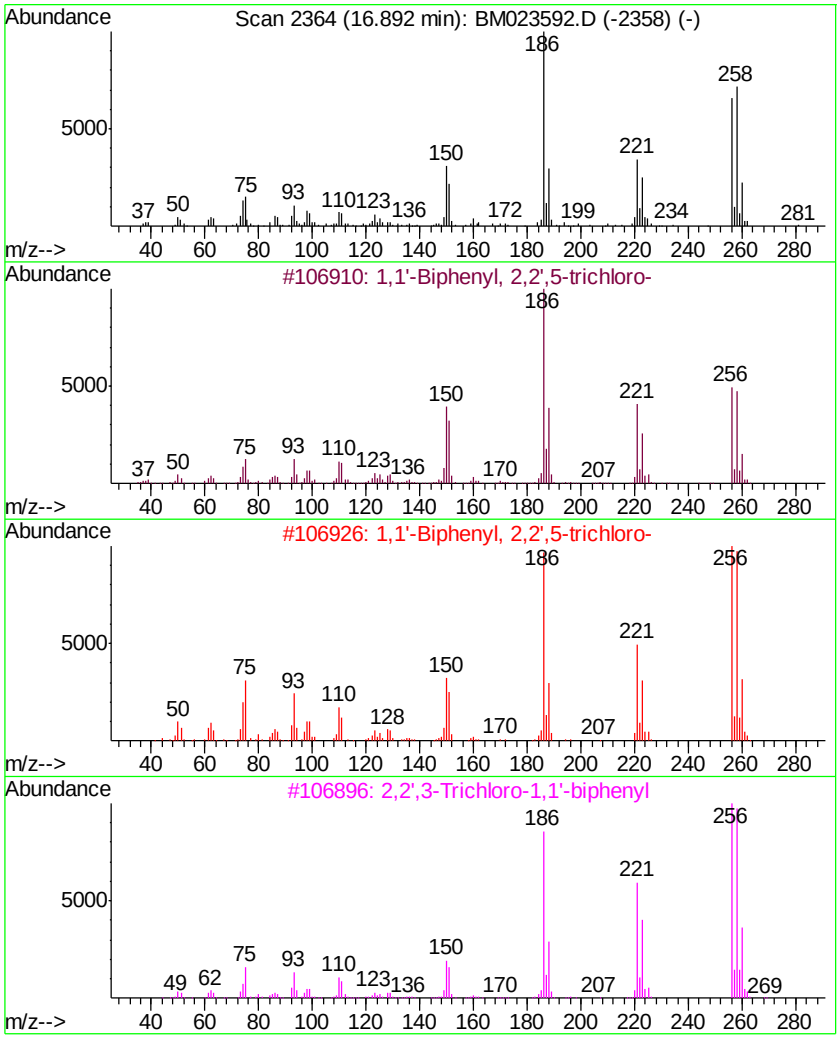
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

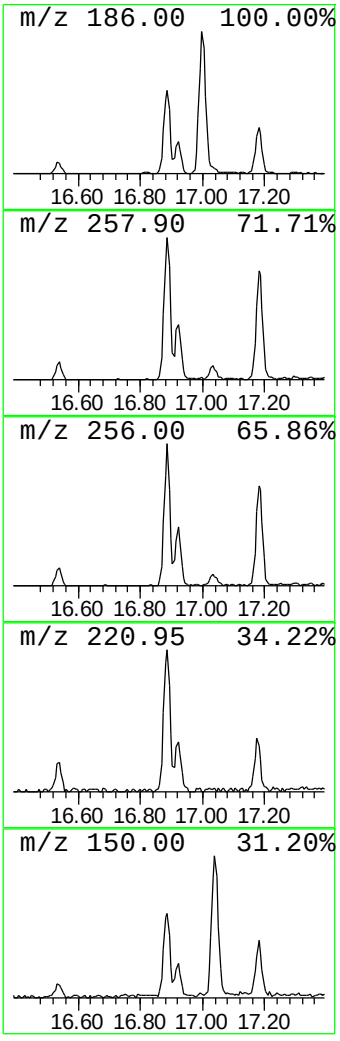
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1,1'-Biphenyl, 2,2',5-trich... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.89	2.04 ng/ul	466041	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99
3	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99
5	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	98





Library Search Compound Report

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Instrument : BNA_M
ClientSampleId : BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

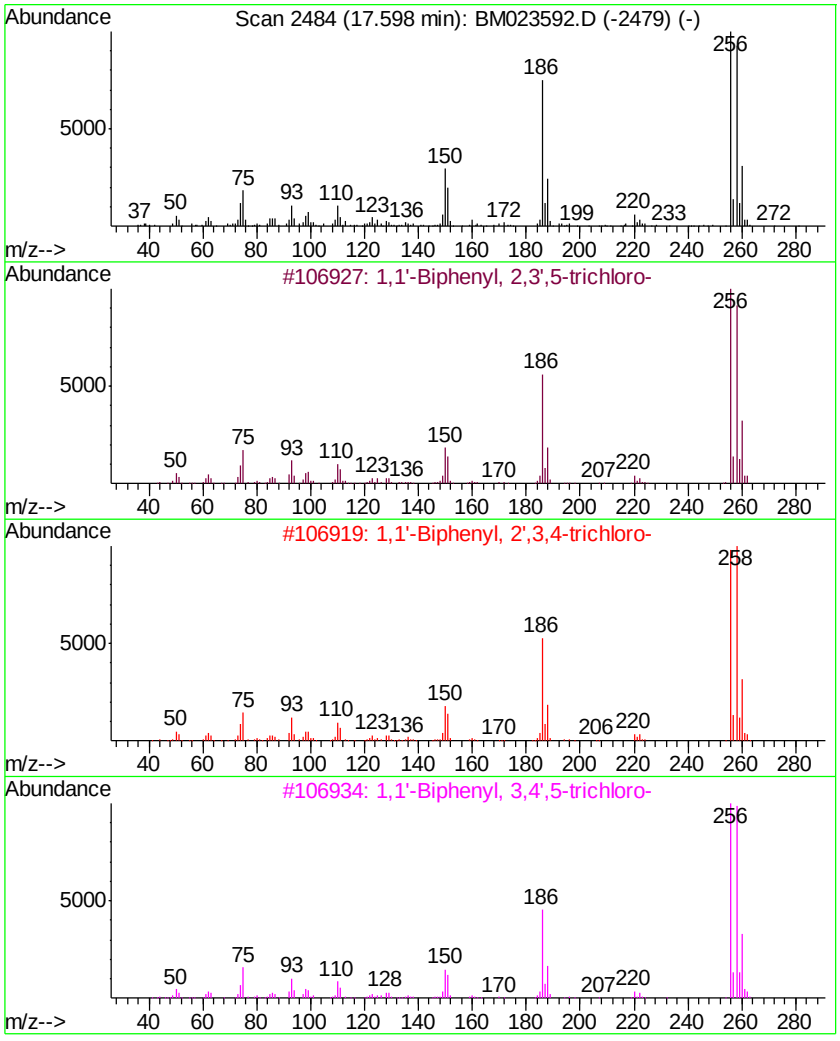
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

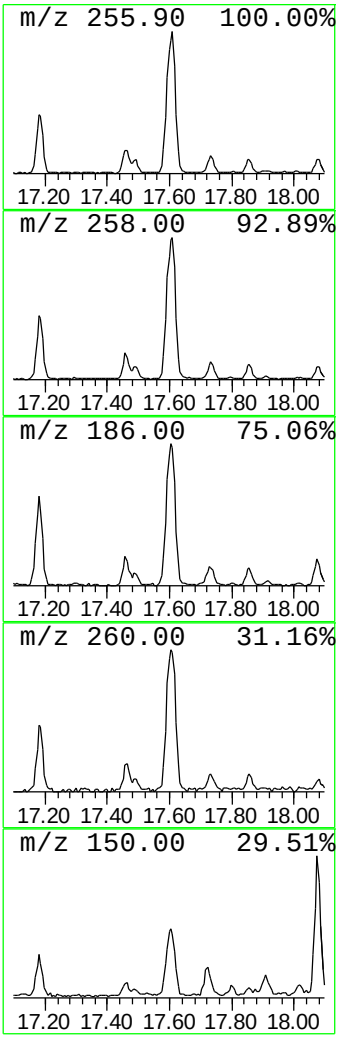
****

Peak Number 3 1,1'-Biphenyl, 2,3',5-trich... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.60	3.48 ng/ul	795458	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99
3	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99
4	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99
5	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	99





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Quant Title : SVOA CALIBRATION

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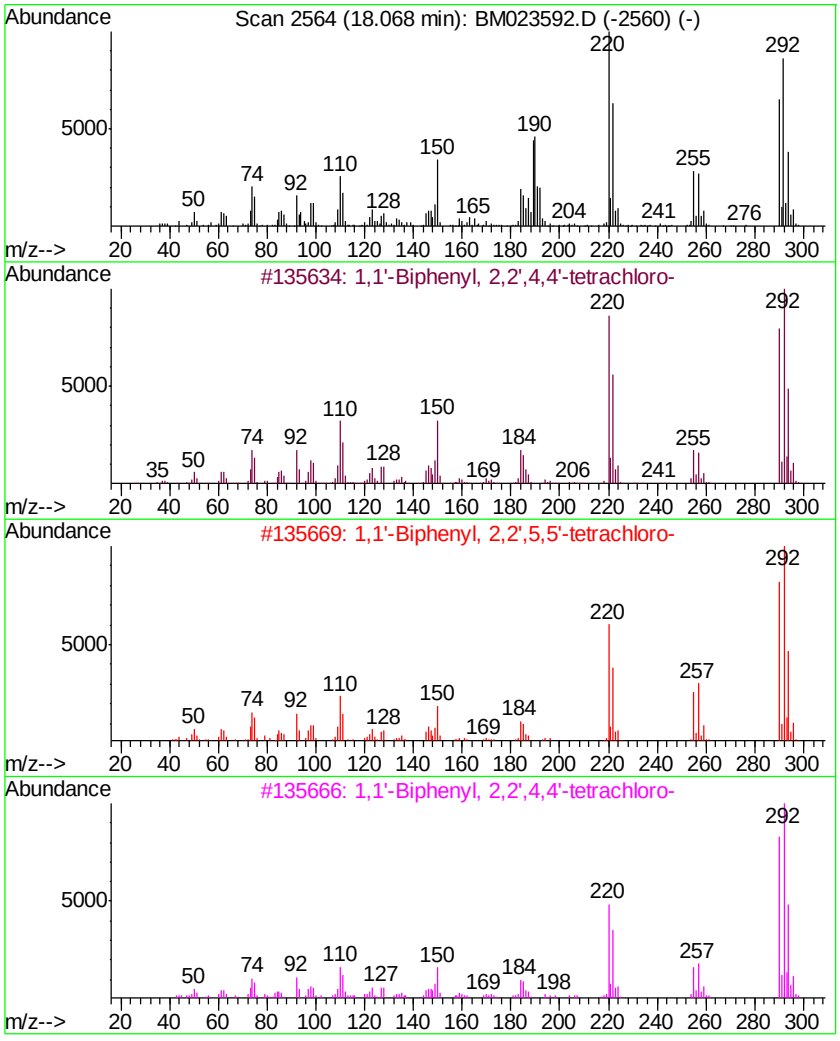
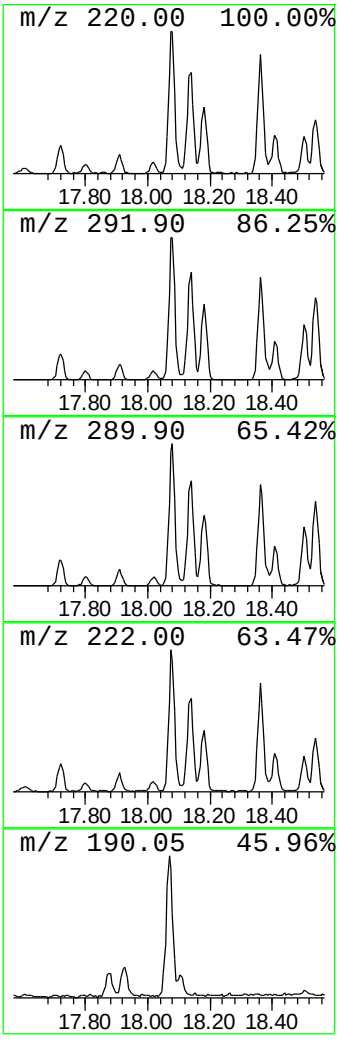
****

Peak Number 4 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	6.45 ng/ul	1473610	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
4	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

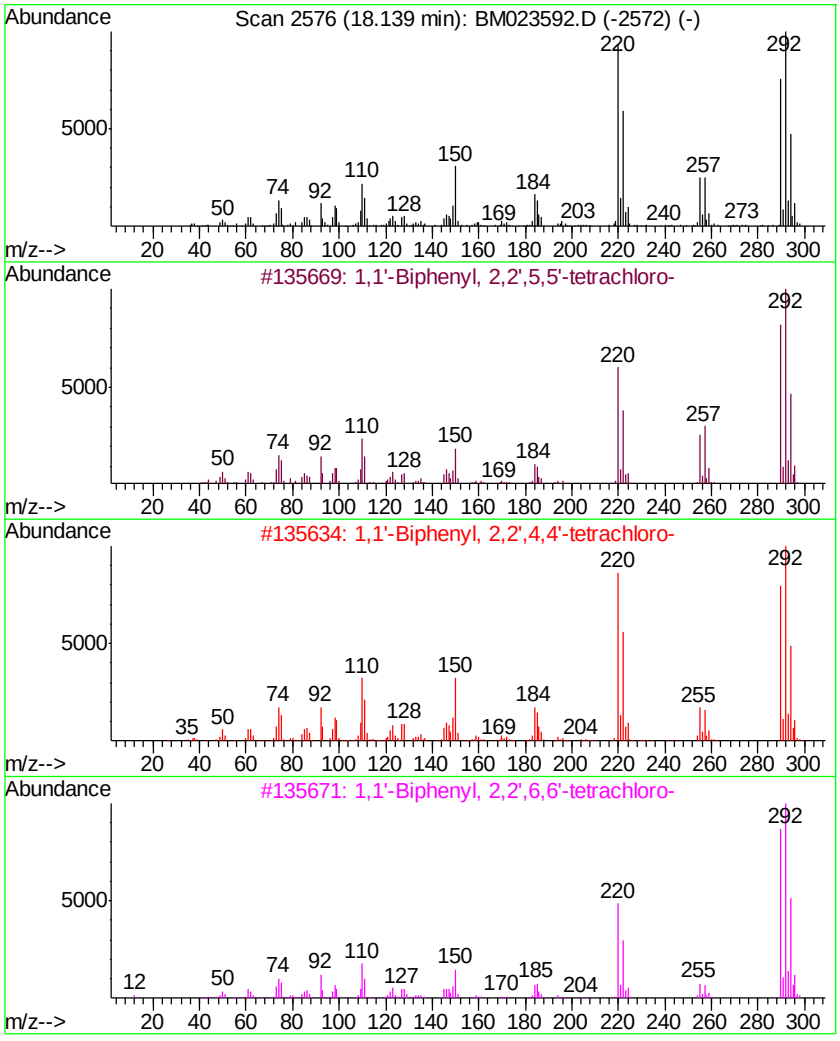
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

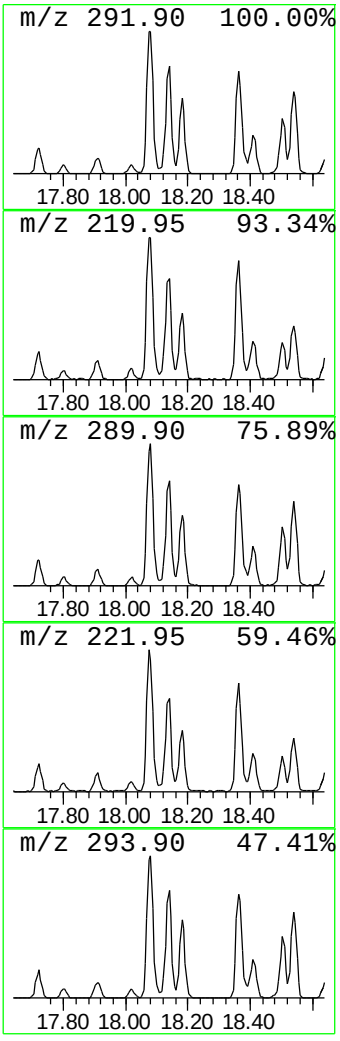
****

Peak Number 5 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.14	3.41 ng/ul	778920	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
4	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
5	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99





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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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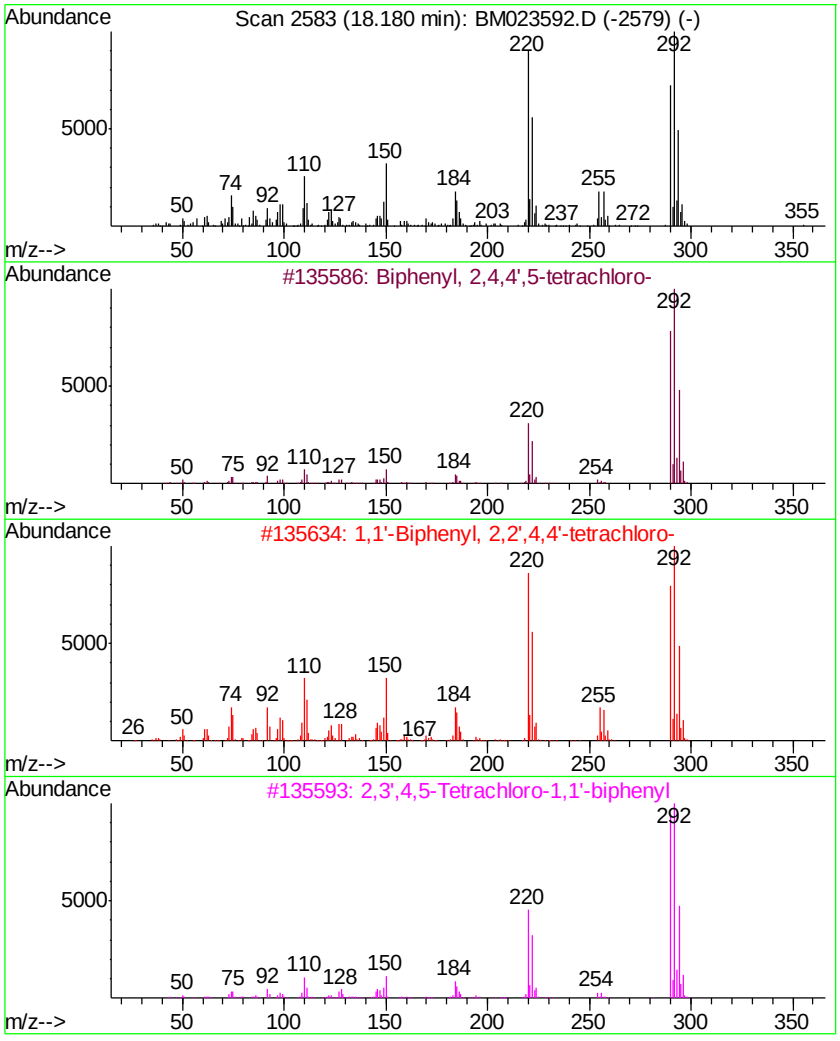
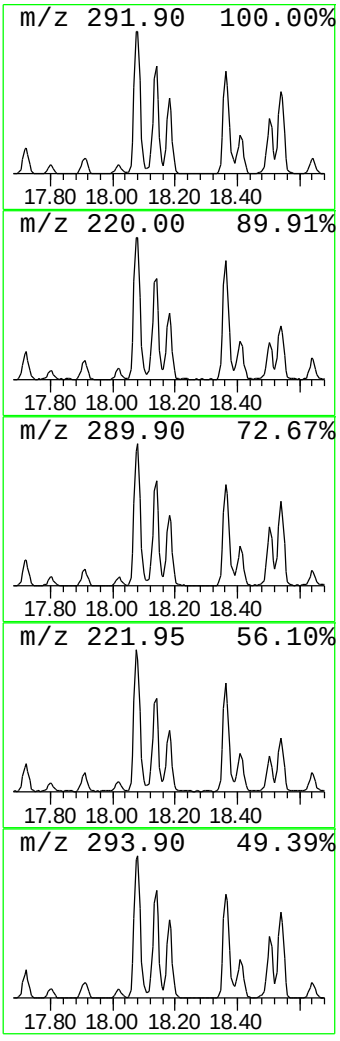
Peak Number 6 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.18	2.42 ng/ul	552781	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
4	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	99
5	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99





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Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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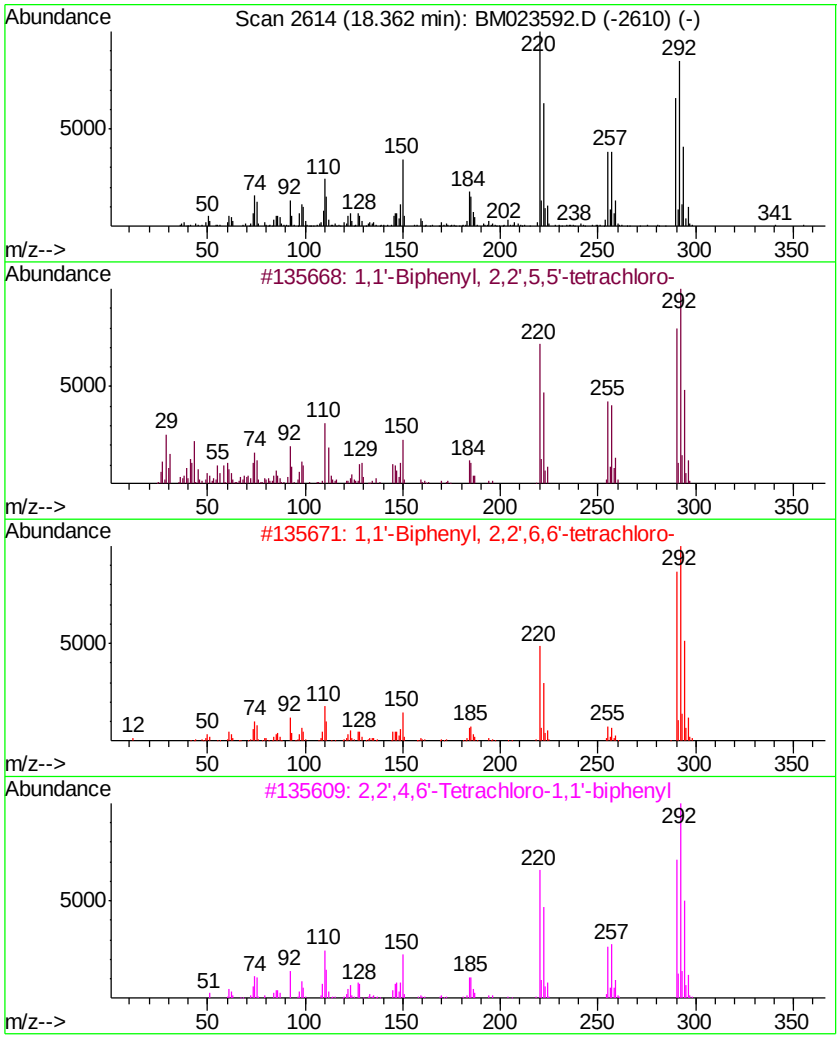
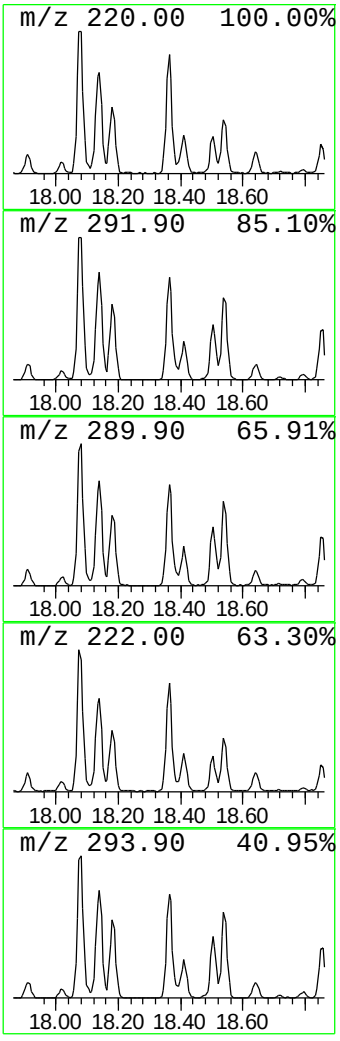
Peak Number 7 1,1'-Biphenyl, 2,2',6,6'-te... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	4.63 ng/ul	1058630	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
4	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110419\
Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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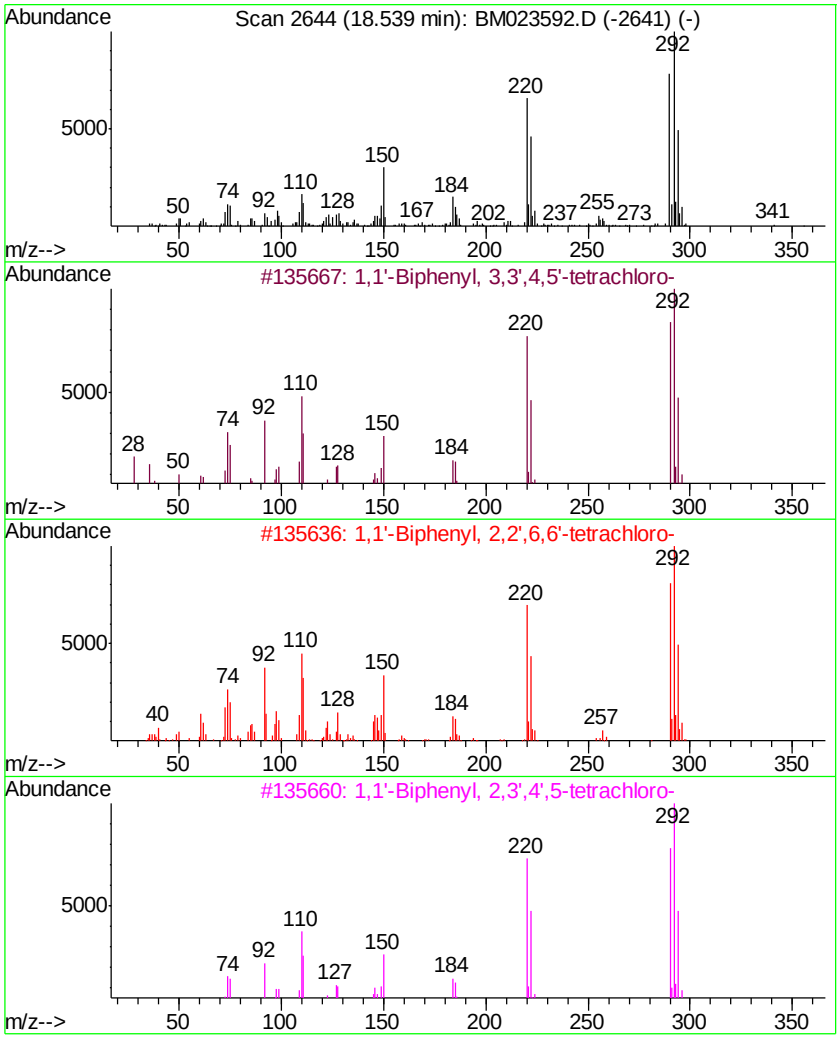
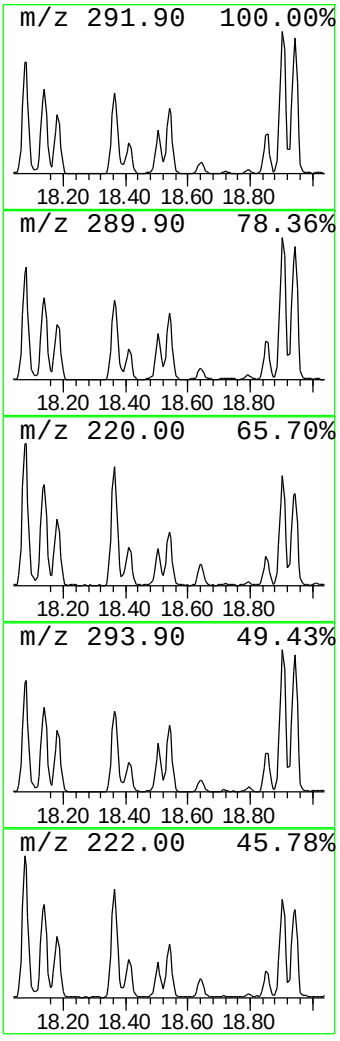
Peak Number 8 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	2.22 ng/ul	508011	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
3	1,1'-Biphenyl, 2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	99
4	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99
5	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110419\
Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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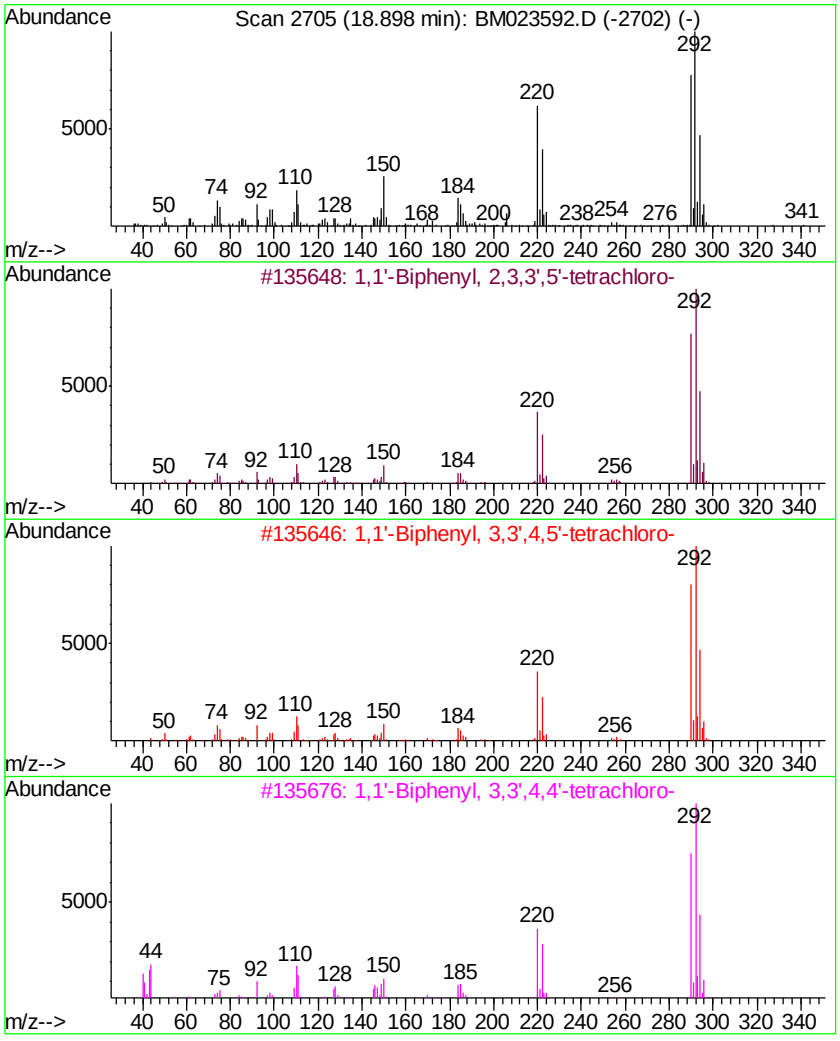
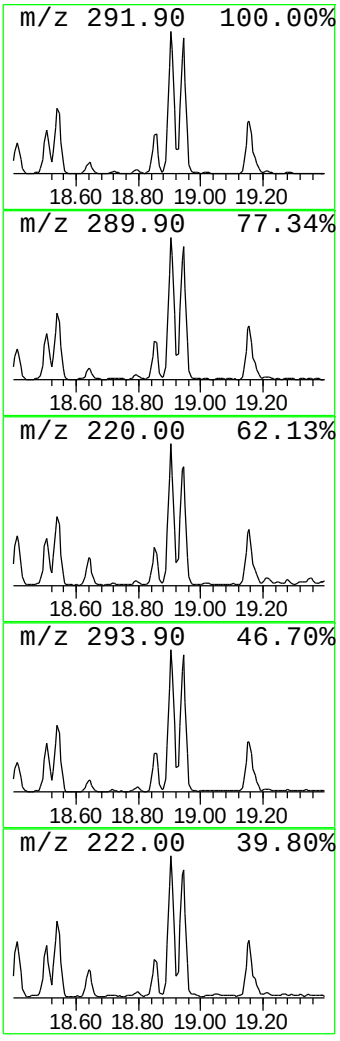
Peak Number 9 1,1'-Biphenyl, 2,3,3',5'-te... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.90	4.84 ng/ul	1107220	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
2	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
3	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99
4	2,3,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	99
5	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	99





Library Search Compound Report

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Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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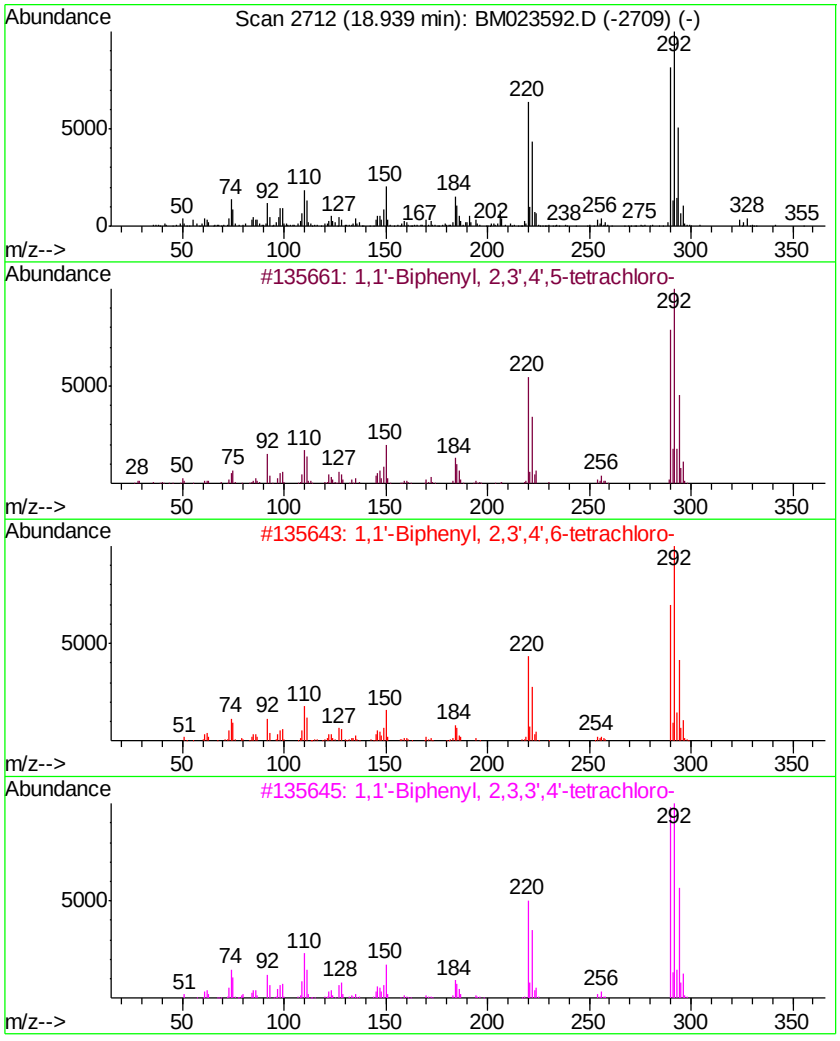
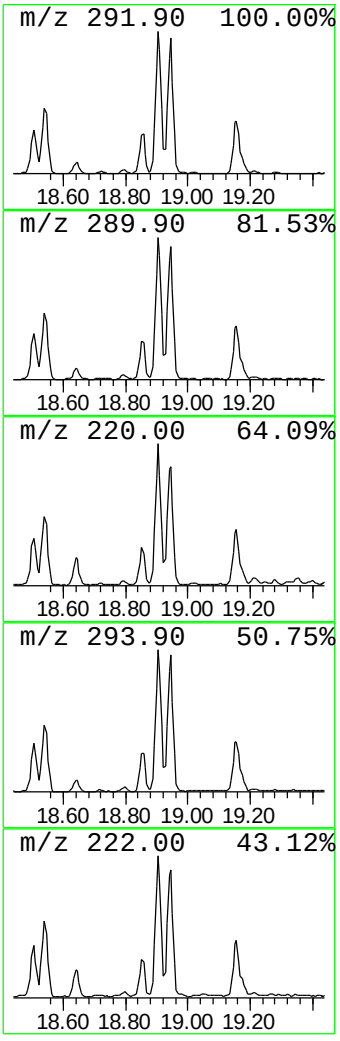
Peak Number 10 1,1'-Biphenyl, 2,3',4',5-te... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.94	5.20 ng/ul	1187690	Phenanthrene-d10	17.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	99
2	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99
3	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	99
4	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99
5	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	99





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM110419\
Data File : BM023592.D
Acq On : 04 Nov 2019 22:57
Operator : JU
Sample : K5403-01
Misc : GCMS CONFIRMATION
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB92

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Str...	3.07	7.4	ng/ul	1057100	1	7.60	2855540	20.0
1,1'-Biphenyl, 2,...	16.89	2.0	ng/ul	466041	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	17.60	3.5	ng/ul	795458	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	18.07	6.5	ng/ul	1473610	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	18.14	3.4	ng/ul	778920	4	17.00	4571730	20.0
Biphenyl, 2,4,4',...	18.18	2.4	ng/ul	552781	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	18.36	4.6	ng/ul	1058630	4	17.00	4571730	20.0
1,1'-Biphenyl, 3,...	18.54	2.2	ng/ul	508011	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	18.90	4.8	ng/ul	1107220	4	17.00	4571730	20.0
1,1'-Biphenyl, 2,...	18.94	5.2	ng/ul	1187690	4	17.00	4571730	20.0