

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050619\
 Data File : BN005486.D
 Acq On : 06 May 2019 10:45
 Operator : JU/SJ
 Sample : K2604-06DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAJ80DL

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.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	5.299	388	393	401	rVB	41795	81875	2.16%	0.176%
2	5.628	444	449	461	rVB4	89951	180824	4.77%	0.389%
3	6.781	638	645	659	rVB	196269	385623	10.17%	0.829%
4	6.940	666	672	680	rBV	154444	240413	6.34%	0.517%
5	7.134	699	705	713	rVB	211780	332916	8.78%	0.716%
6	7.252	719	725	733	rVB2	100641	187404	4.94%	0.403%
7	7.340	733	740	745	rVB3	32474	50296	1.33%	0.108%
8	7.593	776	783	792	rBV	1067747	1697916	44.77%	3.652%
9	7.705	798	802	806	rVB3	16462	27590	0.73%	0.059%
10	8.116	866	872	882	rVB	327164	520447	13.72%	1.119%
11	8.299	897	903	915	rBV	193905	355357	9.37%	0.764%
12	8.681	963	968	971	rBV4	15251	21031	0.55%	0.045%
13	8.734	971	977	986	rVV	187683	342417	9.03%	0.736%
14	8.852	992	997	1006	rVB2	21027	42052	1.11%	0.090%
15	9.169	1046	1051	1053	rBV4	13910	21651	0.57%	0.047%
16	9.263	1062	1067	1071	rBV3	15052	21691	0.57%	0.047%
17	9.340	1075	1080	1084	rBV3	17881	30336	0.80%	0.065%
18	9.452	1092	1099	1106	rBV	144940	250251	6.60%	0.538%
19	9.793	1147	1157	1163	rBV4	67625	159557	4.21%	0.343%
20	9.857	1163	1168	1173	rBV4	53554	106885	2.82%	0.230%
21	9.987	1184	1190	1199	rBV	216657	408544	10.77%	0.879%
22	10.063	1199	1203	1212	rVB5	16583	33286	0.88%	0.072%
23	10.352	1245	1252	1256	rBV	1514797	2519890	66.44%	5.419%
24	10.404	1256	1261	1269	rVV	2311316	3792490	100.00%	8.156%
25	10.499	1270	1277	1290	rVB2	190937	427385	11.27%	0.919%
26	11.187	1389	1394	1401	rBV3	27144	58003	1.53%	0.125%
27	11.293	1405	1412	1416	rVB6	9365	21878	0.58%	0.047%
28	11.399	1421	1430	1434	rBV5	17379	34342	0.91%	0.074%
29	11.681	1475	1478	1484	rVB7	10352	19625	0.52%	0.042%
30	11.798	1493	1498	1505	rVV2	33570	58913	1.55%	0.127%
31	11.869	1507	1510	1514	rVV5	13627	25428	0.67%	0.055%
32	12.022	1527	1536	1547	rBV	1076205	1741998	45.93%	3.746%
33	12.151	1551	1558	1564	rVB5	8759	21607	0.57%	0.046%
34	12.246	1564	1574	1584	rBV	603970	993076	26.19%	2.136%

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Integrator: RTE
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35	12.763	1658	1662	1667	rBV6	12956	20771	0.55%	0.045%
36	13.057	1704	1712	1722	rBV2	174446	298258	7.86%	0.641%
37	13.251	1738	1745	1756	rVB	58243	139543	3.68%	0.300%
38	13.387	1759	1768	1769	rBV	158595	233758	6.16%	0.503%
39	13.463	1777	1781	1786	rVB4	19369	29641	0.78%	0.064%
40	13.545	1786	1795	1800	rBV	264673	466725	12.31%	1.004%
41	13.598	1800	1804	1807	rVV	162644	238494	6.29%	0.513%
42	13.640	1807	1811	1815	rVV	479826	732431	19.31%	1.575%
43	13.681	1815	1818	1829	rVV3	250936	456885	12.05%	0.983%
44	13.781	1829	1835	1839	rVV	125668	207453	5.47%	0.446%
45	13.816	1839	1841	1846	rVV3	39672	47302	1.25%	0.102%
46	13.916	1851	1858	1860	rBV	575806	871820	22.99%	1.875%
47	13.945	1860	1863	1877	rVB	661363	971977	25.63%	2.090%
48	14.145	1892	1897	1901	rBV	35788	48328	1.27%	0.104%
49	14.228	1901	1911	1918	rBV2	2167497	3435286	90.58%	7.388%
50	14.287	1918	1921	1925	rVV2	69492	108959	2.87%	0.234%
51	14.322	1925	1927	1932	rVB2	37003	44204	1.17%	0.095%
52	14.469	1942	1952	1959	rBV6	71410	223719	5.90%	0.481%
53	14.634	1970	1980	1986	rBV	97532	178071	4.70%	0.383%
54	14.710	1988	1993	1997	rBV2	52287	71633	1.89%	0.154%
55	14.839	2009	2015	2021	rBV2	92526	200724	5.29%	0.432%
56	14.910	2023	2027	2030	rVV	40844	57703	1.52%	0.124%
57	14.940	2030	2032	2036	rVB2	23202	26017	0.69%	0.056%
58	15.004	2038	2043	2045	rBV4	33234	57567	1.52%	0.124%
59	15.028	2045	2047	2050	rVB3	16552	17785	0.47%	0.038%
60	15.104	2056	2060	2067	rVB	128669	174306	4.60%	0.375%
61	15.175	2067	2072	2075	rBV5	23159	45746	1.21%	0.098%
62	15.228	2075	2081	2086	rVV	772765	1149102	30.30%	2.471%
63	15.281	2086	2090	2102	rVB	422992	704452	18.57%	1.515%
64	15.375	2102	2106	2109	rBV4	28395	44245	1.17%	0.095%
65	15.410	2109	2112	2115	rVB5	28365	32993	0.87%	0.071%
66	15.492	2121	2126	2129	rBV4	24328	48903	1.29%	0.105%
67	15.581	2136	2141	2146	rBV3	42072	88668	2.34%	0.191%
68	15.698	2155	2161	2164	rBV	108589	168550	4.44%	0.362%
69	15.804	2176	2179	2186	rVB8	14009	22483	0.59%	0.048%
70	15.969	2202	2207	2210	rBV4	63003	93877	2.48%	0.202%
71	16.286	2255	2261	2266	rBV2	80759	168056	4.43%	0.361%

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Integration Parameters: LSCINT.P

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72	16.339	2267	2270	2275	rVB	64926	87304	2.30%	0.188%
73	16.439	2283	2287	2291	rBV	49486	73242	1.93%	0.158%
74	16.645	2319	2322	2327	rVV3	35963	52242	1.38%	0.112%
75	16.722	2331	2335	2339	rVV7	16373	30290	0.80%	0.065%
76	16.763	2339	2342	2344	rVV2	45885	56713	1.50%	0.122%
77	16.798	2344	2348	2360	rVB	134293	240565	6.34%	0.517%
78	16.986	2373	2380	2383	rBV	2589161	3780605	99.69%	8.131%
79	17.028	2383	2387	2391	rVV	1525356	2174792	57.34%	4.677%
80	17.081	2391	2396	2400	rVV2	750673	1215945	32.06%	2.615%
81	17.116	2400	2402	2408	rVV	355787	447940	11.81%	0.963%
82	17.181	2409	2413	2418	rVB5	37325	71509	1.89%	0.154%
83	17.398	2448	2450	2455	rVB5	22108	21920	0.58%	0.047%
84	17.510	2465	2469	2472	rBV3	58777	73772	1.95%	0.159%
85	17.569	2475	2479	2487	rVV2	67071	106841	2.82%	0.230%
86	17.704	2500	2502	2509	rVB2	46425	79087	2.09%	0.170%
87	17.863	2524	2529	2533	rBV	242223	351485	9.27%	0.756%
88	17.910	2533	2537	2542	rVB	289188	453875	11.97%	0.976%
89	17.992	2548	2551	2556	rVB	93485	114361	3.02%	0.246%
90	18.057	2556	2562	2566	rBV2	304133	564515	14.89%	1.214%
91	18.092	2566	2568	2574	rVB	143787	184547	4.87%	0.397%
92	18.369	2612	2615	2619	rBV	101766	124175	3.27%	0.267%
93	18.792	2683	2687	2692	rBV2	111628	191292	5.04%	0.411%
94	19.057	2728	2732	2738	rBV	497788	689514	18.18%	1.483%
95	19.210	2755	2758	2765	rVB	173278	249367	6.58%	0.536%
96	19.398	2785	2790	2792	rBV	768126	1076135	28.38%	2.314%
97	19.427	2792	2795	2801	rVB	641390	882311	23.26%	1.898%
98	19.927	2878	2880	2884	rVB	120387	142628	3.76%	0.307%
99	21.216	3094	3099	3104	rBV	2085302	3078965	81.19%	6.622%
100	23.427	3470	3475	3485	rBV	1377587	2741588	72.29%	5.896%

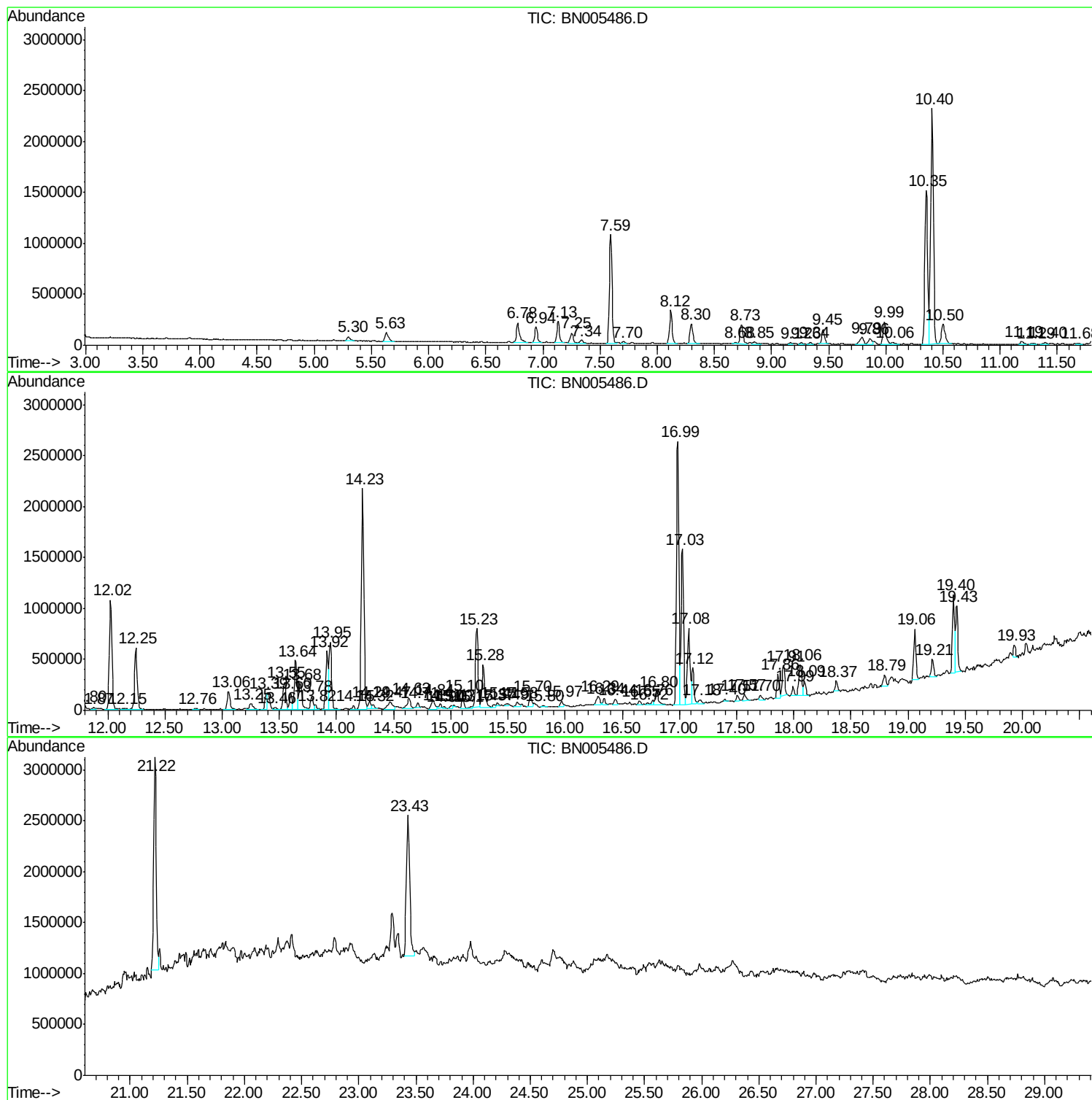
Sum of corrected areas: 46496957

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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampled :
GAJ80DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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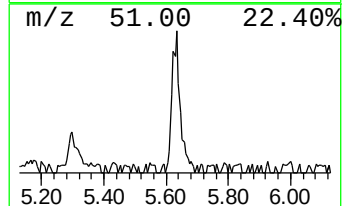
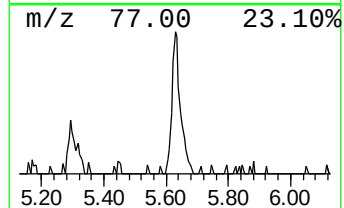
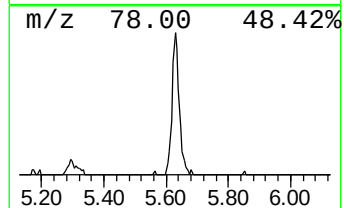
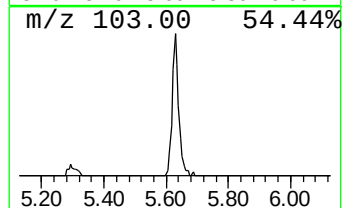
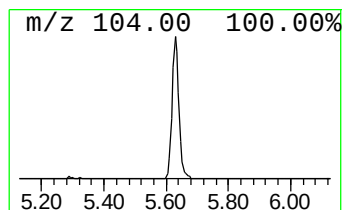
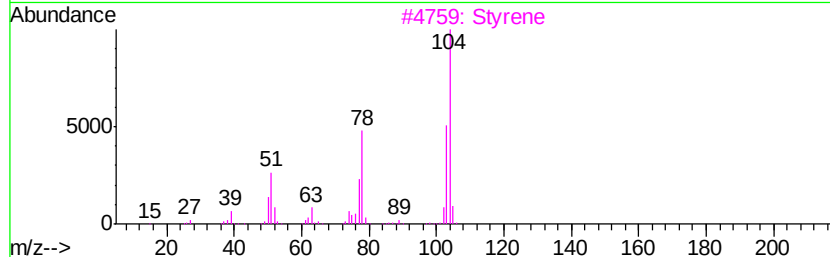
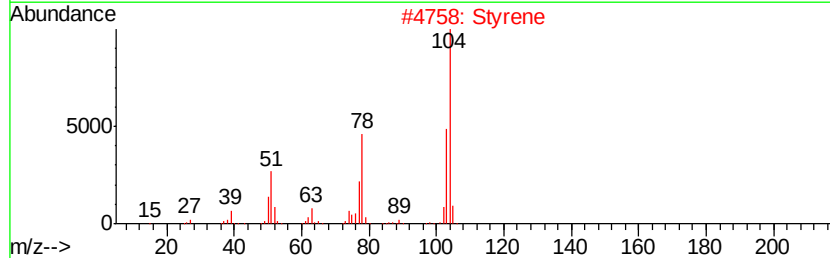
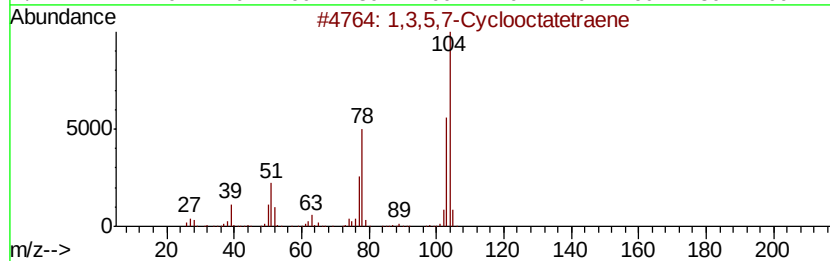
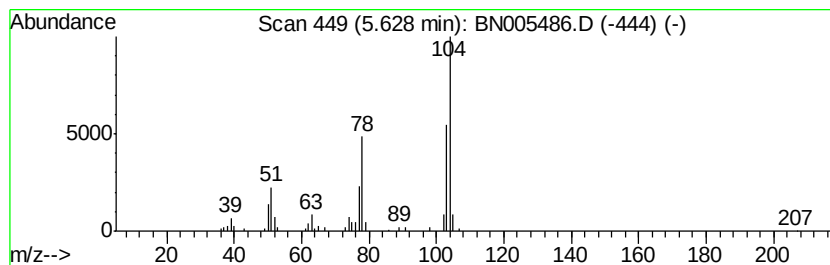
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3,5,7-Cyclooctatetraene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	2.13 ng/ul	180824	1,4-Dichlorobenzene-d4	7.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	97
2			Styrene	104	C8H8	000100-42-5	96
3			Styrene	104	C8H8	000100-42-5	96
4			Styrene	104	C8H8	000100-42-5	95
5			Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	95



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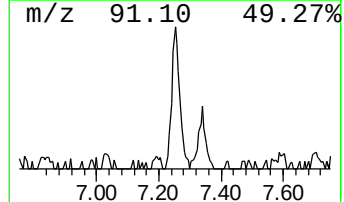
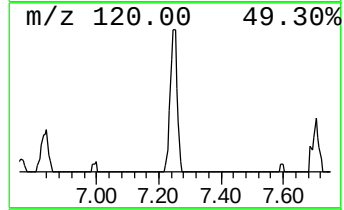
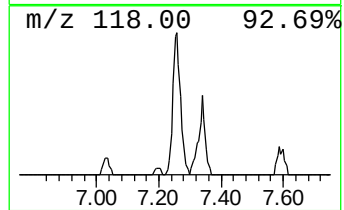
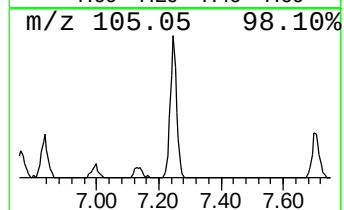
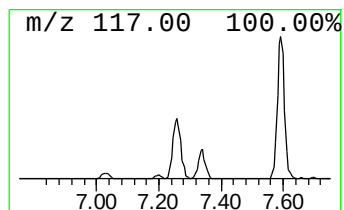
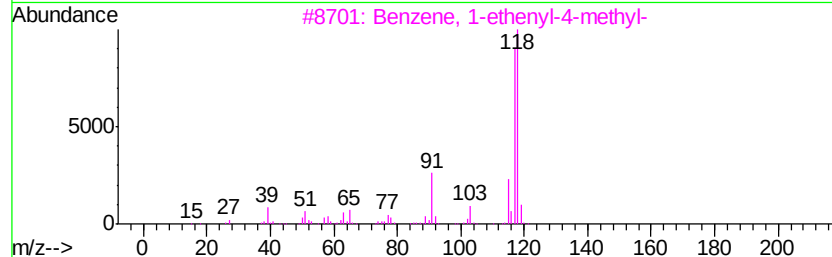
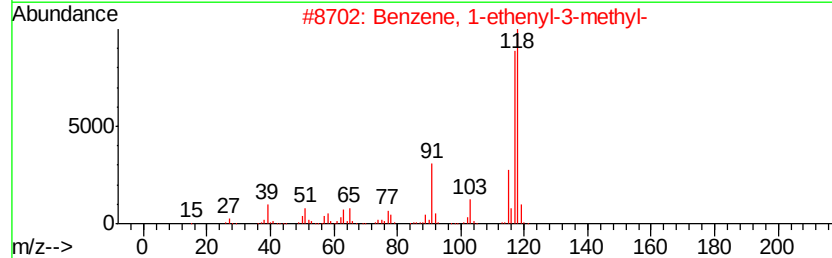
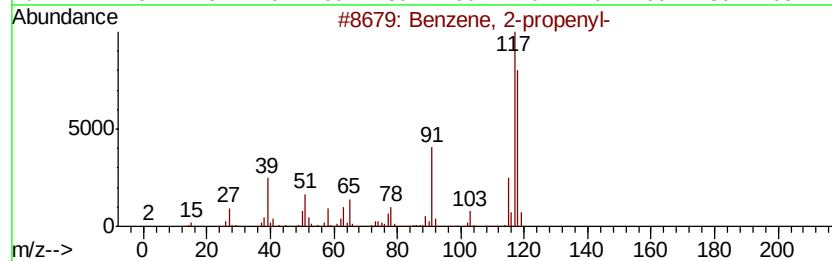
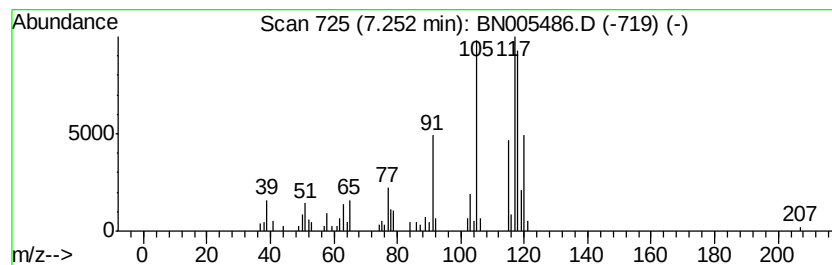
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 2-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	2.21 ng/ul	187404	1,4-Dichlorobenzene-d4	7.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	86
2		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	60
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	60
4		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	60
5		7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	60



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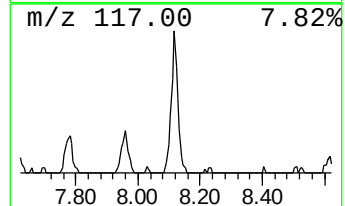
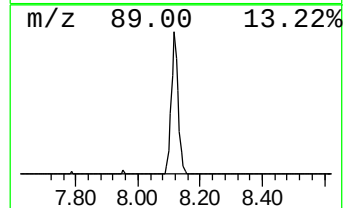
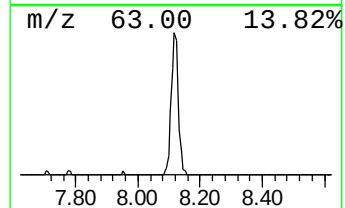
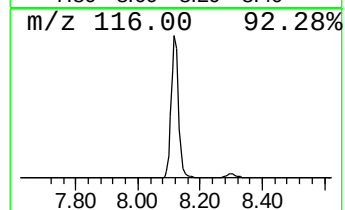
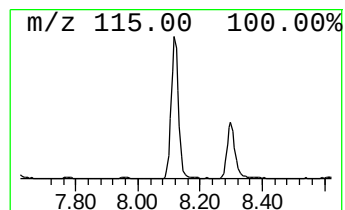
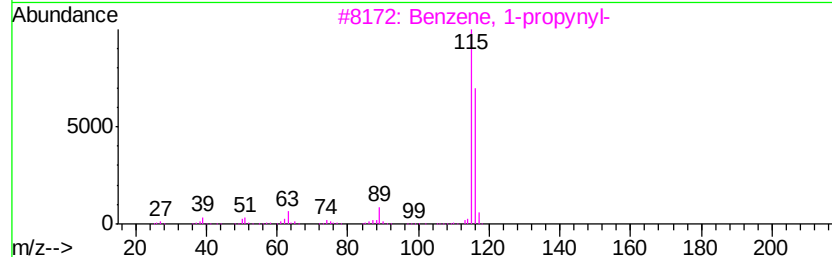
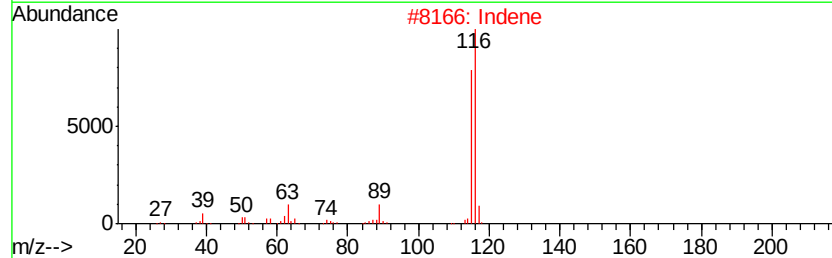
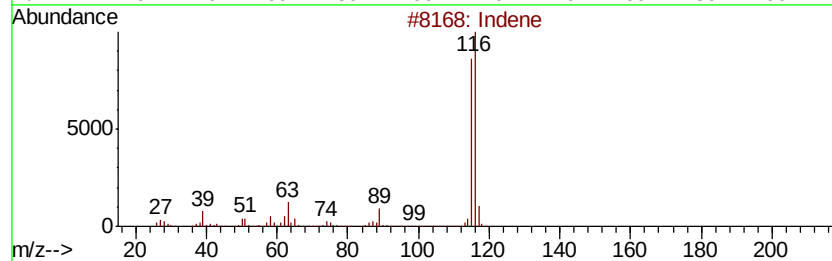
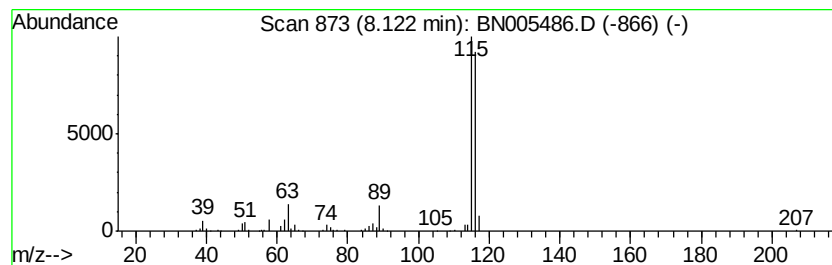
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	6.13 ng/ul	520447	1,4-Dichlorobenzene-d4	7.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene		116	C9H8	000095-13-6	97
2	Indene		116	C9H8	000095-13-6	95
3	Benzene, 1-propynyl-		116	C9H8	000673-32-5	94
4	Benzene, 1-ethynyl-2-methyl-		116	C9H8	000766-47-2	91
5	Benzene, 1-propynyl-		116	C9H8	000673-32-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050619\
Data File : BN005486.D
Acq On : 06 May 2019 10:45
Operator : JU/SJ
Sample : K2604-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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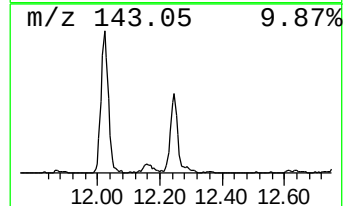
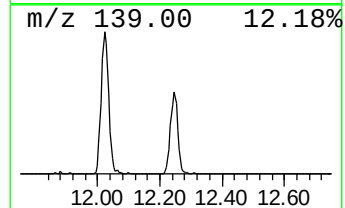
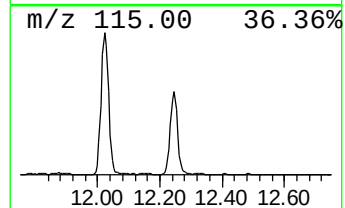
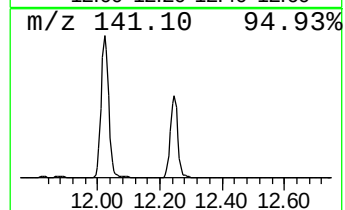
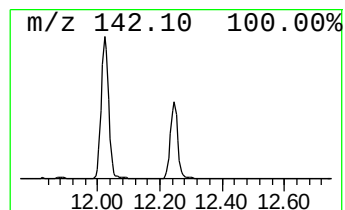
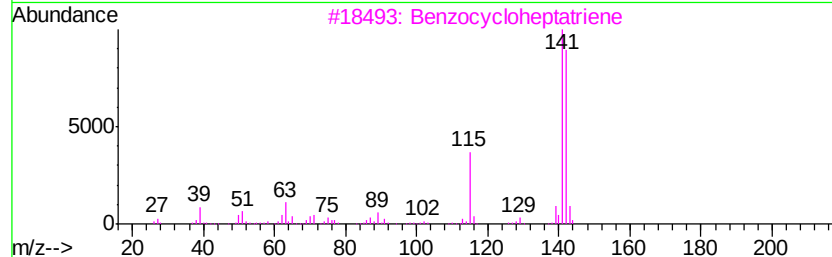
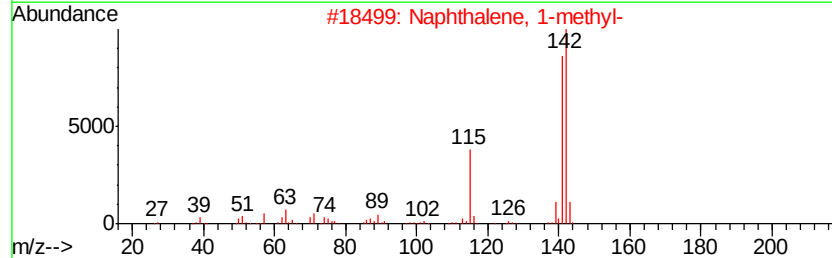
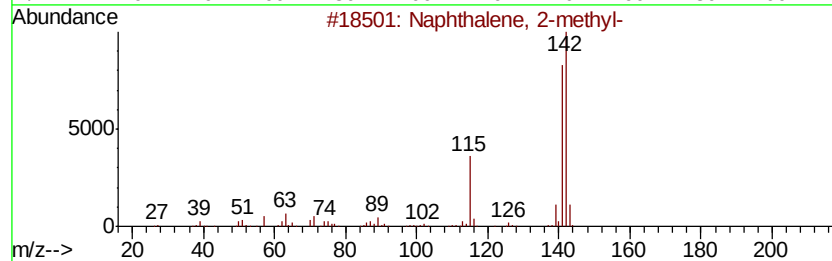
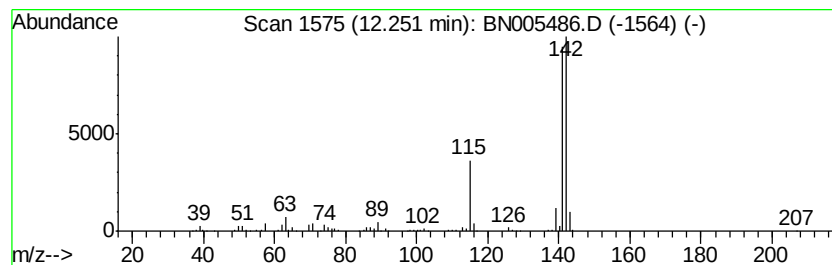
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzocycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	7.88 ng/ul	993076	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050619\
Data File : BN005486.D
Acq On : 06 May 2019 10:45
Operator : JU/SJ
Sample : K2604-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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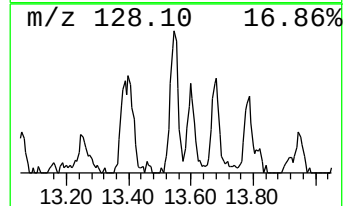
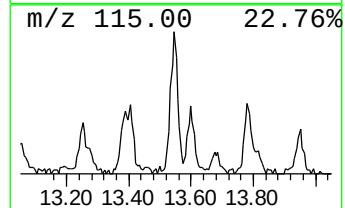
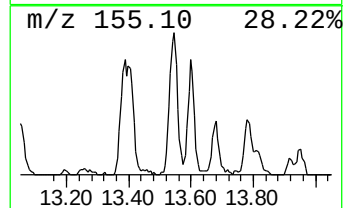
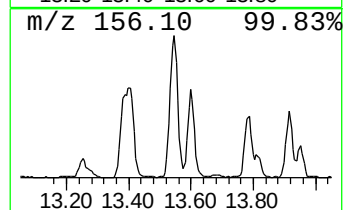
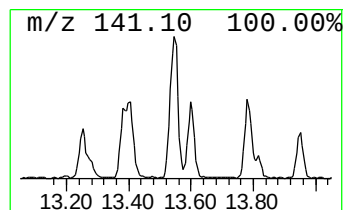
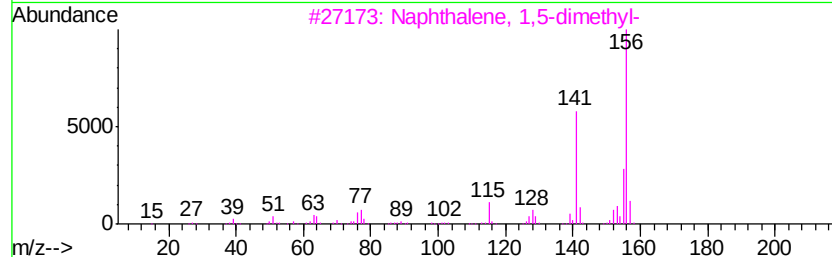
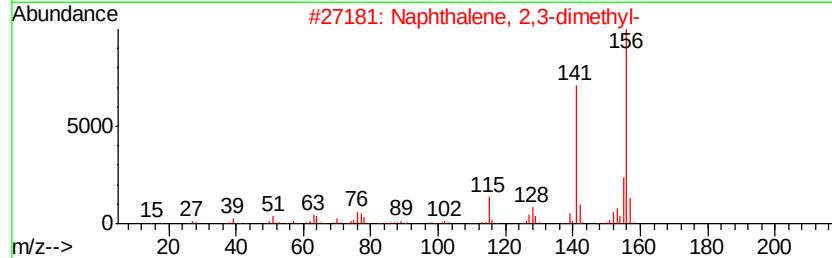
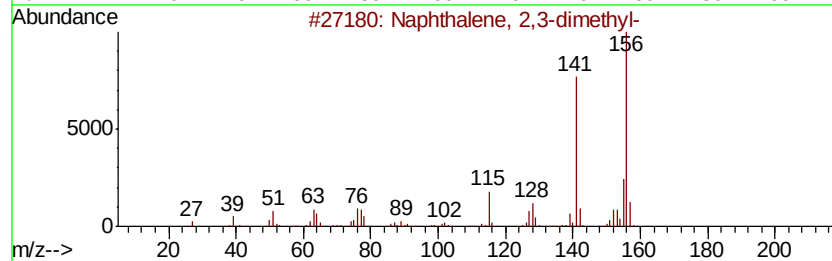
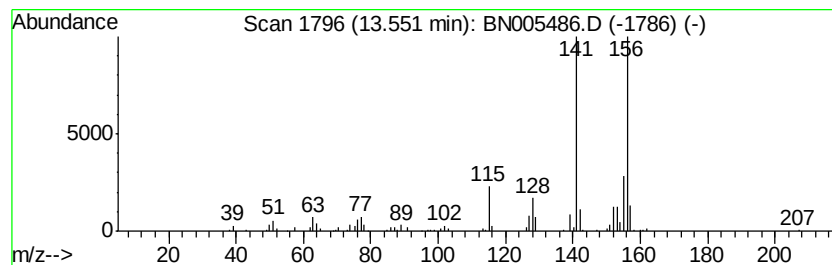
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.55	2.72 ng/ul	466725	Acenaphthene-d10	14.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050619\
Data File : BN005486.D
Acq On : 06 May 2019 10:45
Operator : JU/SJ
Sample : K2604-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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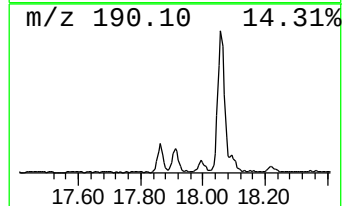
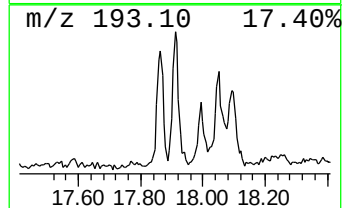
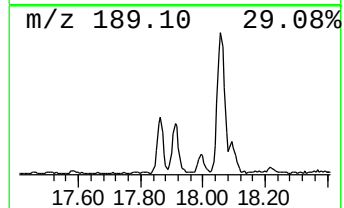
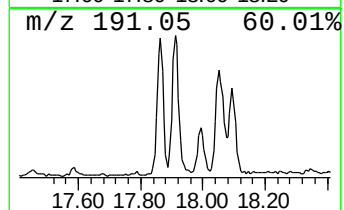
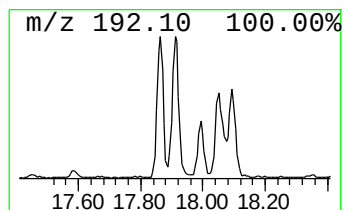
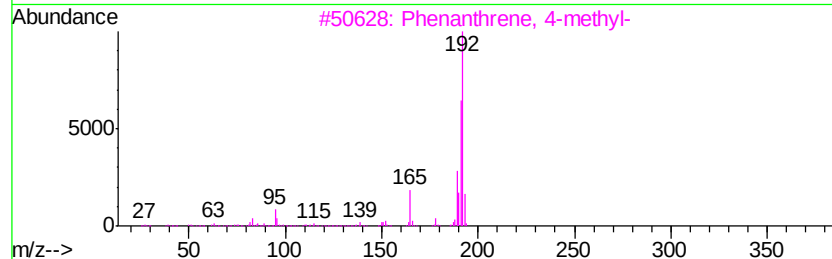
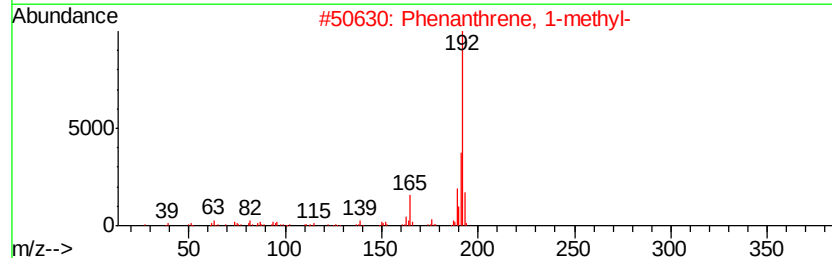
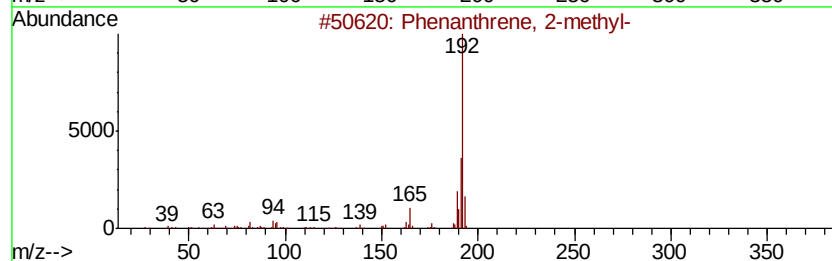
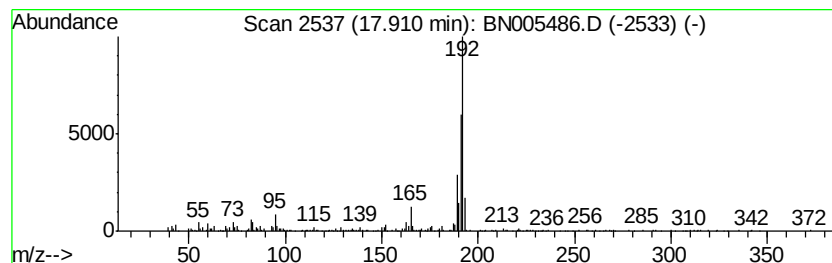
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	2.40 ng/ul	453875	Phenanthrene-d10	16.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050619\
Data File : BN005486.D
Acq On : 06 May 2019 10:45
Operator : JU/SJ
Sample : K2604-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

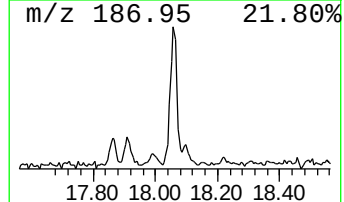
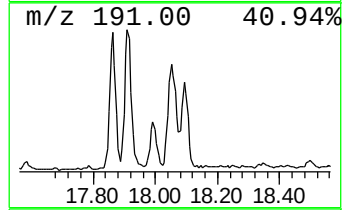
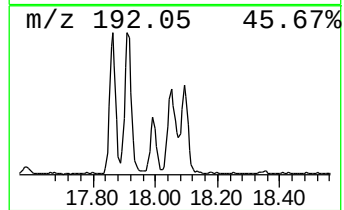
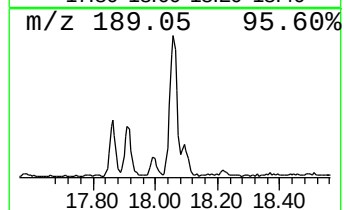
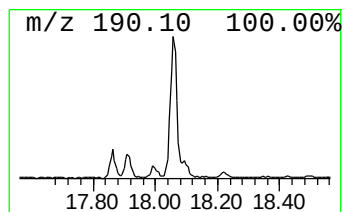
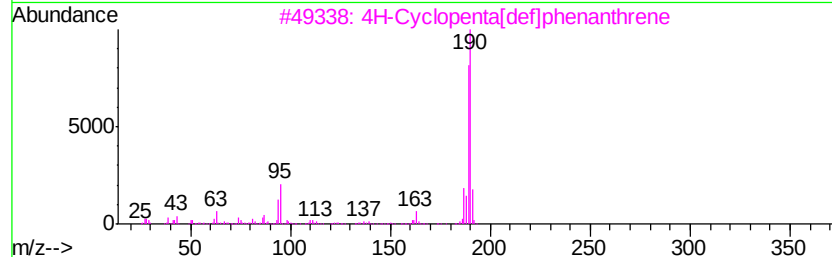
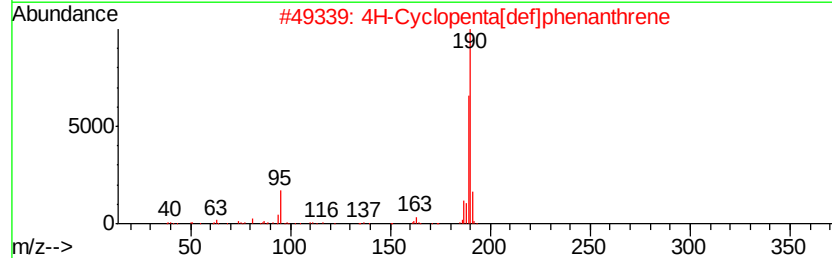
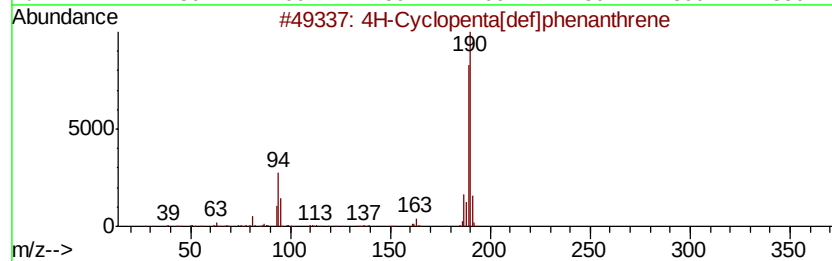
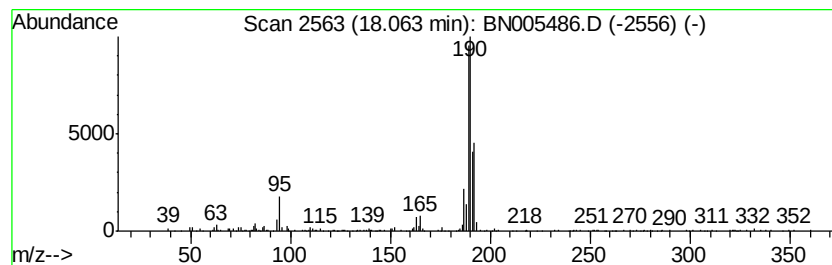
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.		
18.06	2.99 ng/ul	564515	Phenanthrene-d10		16.99		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	86
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4			Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
5			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN050619\
Data File : BN005486.D
Acq On : 06 May 2019 10:45
Operator : JU/SJ
Sample : K2604-06DL 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 2-propenyl-	7.25	2.2	ng/ul	187404	1	7.59	1697920	20.0
Indene	8.12	6.1	ng/ul	520447	1	7.59	1697920	20.0
Benzocycloheptatr...	12.25	7.9	ng/ul	993076	2	10.35	2519890	20.0
Naphthalene, 2,3-...	13.55	2.7	ng/ul	466725	3	14.23	3435290	20.0
Phenanthrene, 2-m...	17.91	2.4	ng/ul	453875	4	16.99	3780610	20.0
4H-Cyclopenta[def...	18.06	3.0	ng/ul	564515	4	16.99	3780610	20.0