

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
 Data File : BN002679.D
 Acq On : 13 Sep 2018 17:17
 Operator : JU/SJ
 Sample : J4864-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0D11

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BN081318MA.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	3.205	32	37	40	rBV	74251	113999	2.77%	0.204%
2	3.234	40	42	47	rVB	43514	53112	1.29%	0.095%
3	6.851	644	657	670	rBV2	235258	528963	12.84%	0.945%
4	7.004	676	683	693	rBV	749306	1148099	27.86%	2.051%
5	7.198	710	716	732	rVB	806219	1325549	32.17%	2.368%
6	7.657	788	794	802	rBV	728234	1196526	29.03%	2.138%
7	7.910	829	837	845	rBV	137702	271077	6.58%	0.484%
8	8.375	905	916	925	rBV	650962	1097039	26.62%	1.960%
9	8.810	983	990	999	rBV	938912	1566518	38.01%	2.799%
10	9.269	1064	1068	1076	rVB	31849	52045	1.26%	0.093%
11	9.522	1104	1111	1123	rBV	796436	1361844	33.05%	2.433%
12	9.628	1123	1129	1137	rBV3	26151	64221	1.56%	0.115%
13	9.840	1161	1165	1173	rVB4	28407	53644	1.30%	0.096%
14	10.063	1197	1203	1211	rBV	1134044	2007108	48.70%	3.586%
15	10.428	1255	1265	1271	rBV	1115016	1918203	46.55%	3.427%
16	10.481	1271	1274	1281	rVB3	76957	127742	3.10%	0.228%
17	10.581	1286	1291	1299	rVB2	75314	155481	3.77%	0.278%
18	10.892	1339	1344	1353	rVB3	96751	176347	4.28%	0.315%
19	11.004	1356	1363	1372	rVV	77027	139616	3.39%	0.249%
20	11.334	1414	1419	1425	rVB5	19469	44781	1.09%	0.080%
21	11.539	1449	1454	1459	rVB2	57950	102900	2.50%	0.184%
22	11.757	1485	1491	1497	rBV	51506	87930	2.13%	0.157%
23	11.963	1522	1526	1533	rBV8	22923	55562	1.35%	0.099%
24	12.063	1539	1543	1554	rVB3	40324	89000	2.16%	0.159%
25	12.228	1563	1571	1579	rBV	216821	397277	9.64%	0.710%
26	12.304	1579	1584	1588	rBV6	40425	75175	1.82%	0.134%
27	12.357	1589	1593	1597	rVV2	55188	84110	2.04%	0.150%
28	12.434	1603	1606	1612	rVB2	40327	53629	1.30%	0.096%
29	12.628	1633	1639	1642	rBV4	31297	54763	1.33%	0.098%
30	12.669	1642	1646	1650	rVB	40027	51402	1.25%	0.092%
31	12.757	1657	1661	1667	rVB3	20480	43397	1.05%	0.078%
32	12.828	1669	1673	1677	rVV3	40455	69709	1.69%	0.125%
33	13.139	1720	1726	1732	rVB3	34039	65451	1.59%	0.117%
34	13.222	1735	1740	1746	rVB	74538	128239	3.11%	0.229%

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35	13.286	1746	1751	1756	rVB4	49155	78192	1.90%	0.140%
36	13.345	1756	1761	1766	rVB	87222	125247	3.04%	0.224%
37	13.710	1817	1823	1828	rBV	1989383	2787272	67.64%	4.979%
38	13.757	1828	1831	1834	rVB	103302	120460	2.92%	0.215%
39	13.845	1842	1846	1849	rBV	98018	161183	3.91%	0.288%
40	13.880	1849	1852	1857	rVB	126506	176257	4.28%	0.315%
41	13.986	1863	1870	1879	rBV	2387754	3581177	86.90%	6.398%
42	14.292	1915	1922	1929	rBV2	1594611	2504103	60.76%	4.474%
43	14.357	1929	1933	1936	rBV	68261	93428	2.27%	0.167%
44	14.545	1957	1965	1972	rBV2	107276	233560	5.67%	0.417%
45	14.828	2010	2013	2021	rVB9	40036	78484	1.90%	0.140%
46	14.916	2022	2028	2035	rBV2	169011	332677	8.07%	0.594%
47	15.045	2046	2050	2053	rBV	90773	149174	3.62%	0.267%
48	15.127	2060	2064	2071	rVB8	47191	89593	2.17%	0.160%
49	15.292	2086	2092	2097	rVV	2914092	4121014	100.00%	7.362%
50	15.339	2097	2100	2105	rVB3	160782	247668	6.01%	0.442%
51	15.433	2111	2116	2119	rBV	520375	739839	17.95%	1.322%
52	15.469	2119	2122	2126	rVV	216400	311115	7.55%	0.556%
53	15.510	2126	2129	2135	rVV5	68926	162948	3.95%	0.291%
54	15.563	2135	2138	2142	rVB	72796	97337	2.36%	0.174%
55	15.645	2148	2152	2153	rBV3	46357	62732	1.52%	0.112%
56	15.857	2186	2188	2200	rVB8	43247	109350	2.65%	0.195%
57	15.992	2208	2211	2215	rBV	73719	94467	2.29%	0.169%
58	16.163	2237	2240	2243	rVB	37488	44234	1.07%	0.079%
59	16.404	2276	2281	2286	rBV3	138151	227535	5.52%	0.406%
60	16.557	2303	2307	2310	rBV4	36099	61142	1.48%	0.109%
61	16.598	2311	2314	2316	rBV3	35256	44098	1.07%	0.079%
62	16.857	2355	2358	2366	rVB4	74905	139428	3.38%	0.249%
63	17.045	2384	2390	2396	rBV	1658525	2382303	57.81%	4.256%
64	17.145	2400	2407	2417	rVB	2782996	3901439	94.67%	6.970%
65	17.951	2539	2544	2553	rBV2	284605	449243	10.90%	0.803%
66	19.239	2759	2763	2773	rBV2	76898	139349	3.38%	0.249%
67	19.445	2792	2798	2805	rBV	2741549	3735468	90.64%	6.673%
68	19.998	2889	2892	2897	rVB	94505	109824	2.66%	0.196%
69	20.380	2954	2957	2962	rVB	76706	88198	2.14%	0.158%
70	20.498	2974	2977	2983	rVB	162116	194510	4.72%	0.347%
71	20.968	3053	3057	3064	rBV	301085	438817	10.65%	0.784%

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72	21.245	3099	3104	3109	rBV	1648653	2115053	51.32%	3.779%
73	21.404	3127	3131	3136	rVB	371064	400201	9.71%	0.715%
74	21.833	3200	3204	3209	rVB	387540	443000	10.75%	0.791%
75	22.292	3277	3282	3290	rBV	789897	1108961	26.91%	1.981%
76	22.415	3299	3303	3308	rVB	418777	552468	13.41%	0.987%
77	22.792	3362	3367	3376	rVB	448032	664804	16.13%	1.188%
78	23.327	3451	3458	3467	rVB2	1943112	4104451	99.60%	7.333%
79	23.468	3475	3482	3492	rVB	1125010	2156526	52.33%	3.853%
80	23.974	3563	3568	3576	rVB	384690	715100	17.35%	1.278%
81	24.703	3687	3692	3700	rVB	270588	541089	13.13%	0.967%

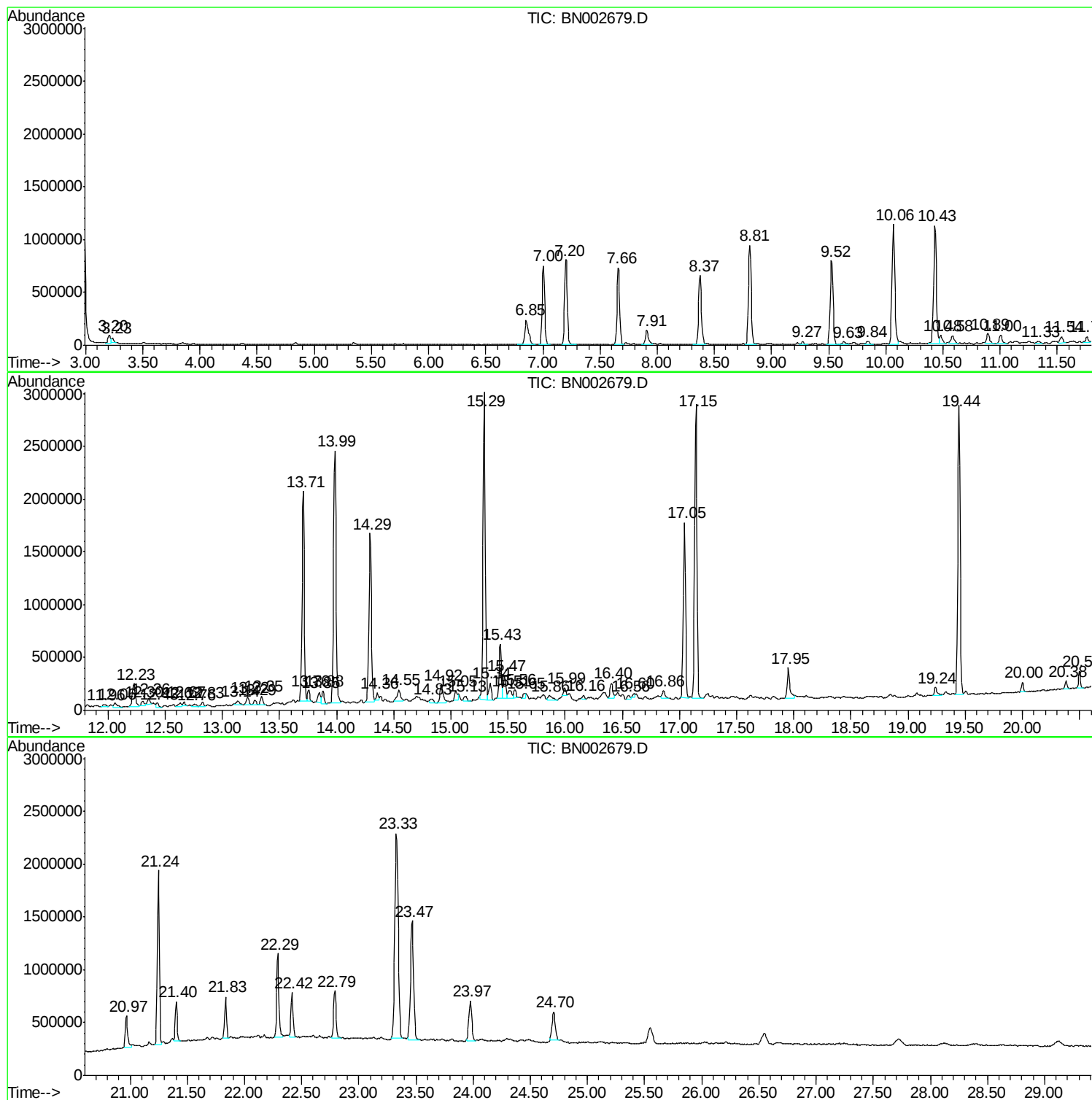
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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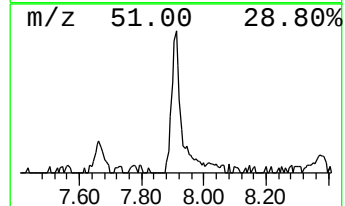
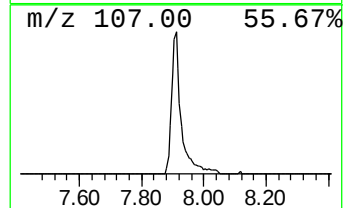
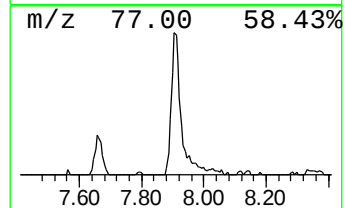
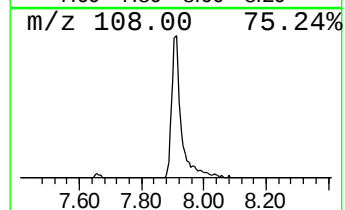
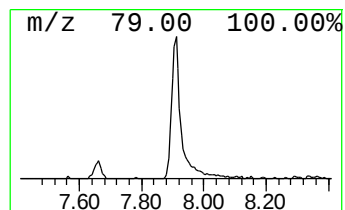
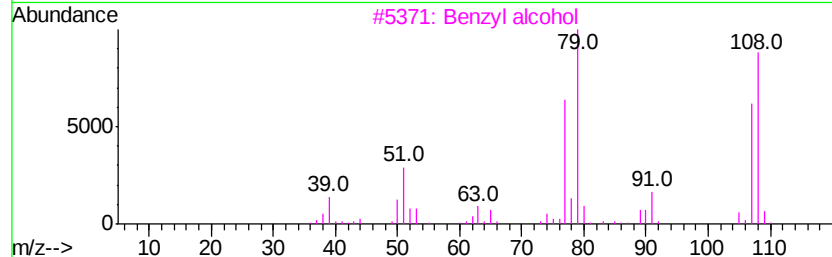
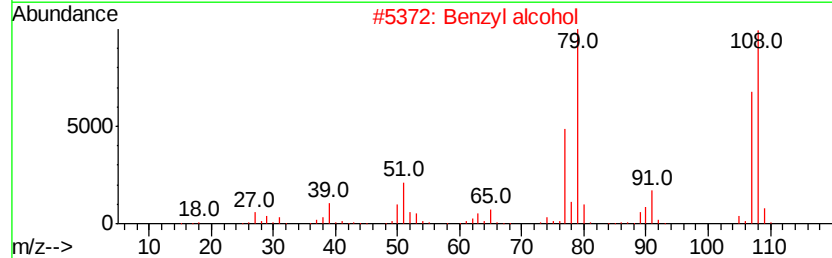
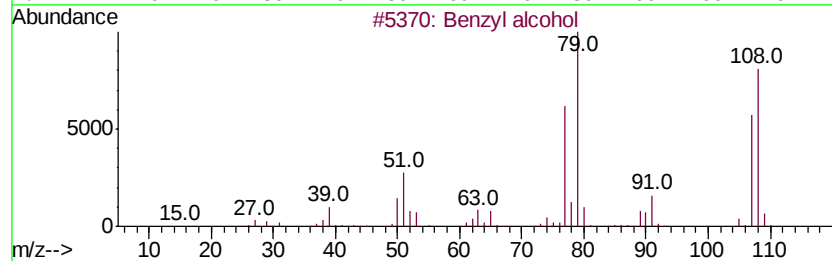
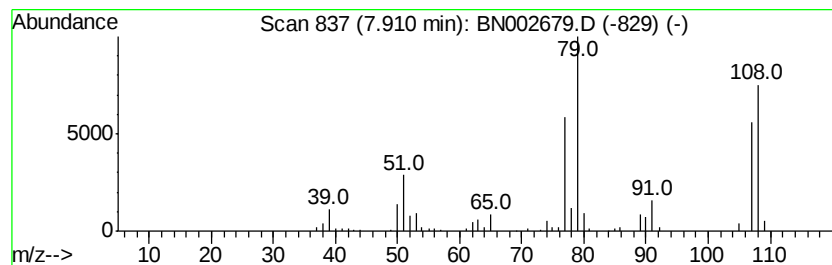
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzyl alcohol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.91	4.53 ng/ul	271077	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	95
2			Benzyl alcohol	108	C7H8O	000100-51-6	94
3			Benzyl alcohol	108	C7H8O	000100-51-6	94
4			Benzyl alcohol	108	C7H8O	000100-51-6	94
5			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



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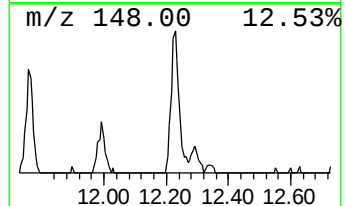
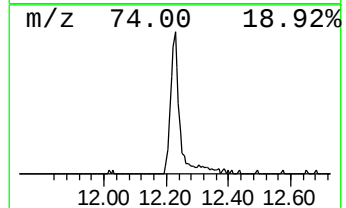
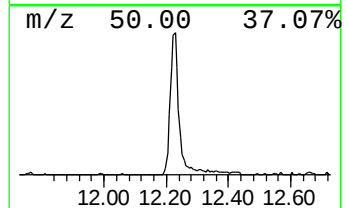
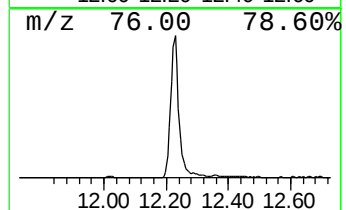
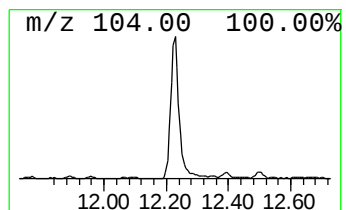
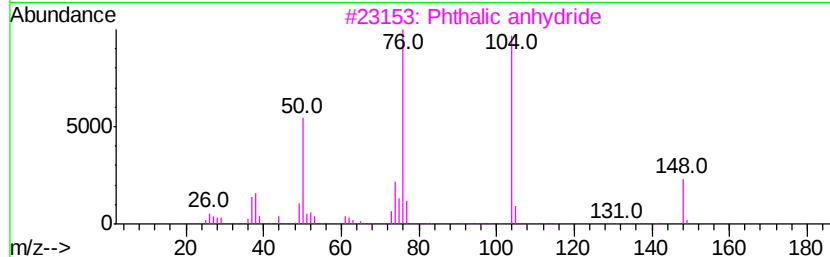
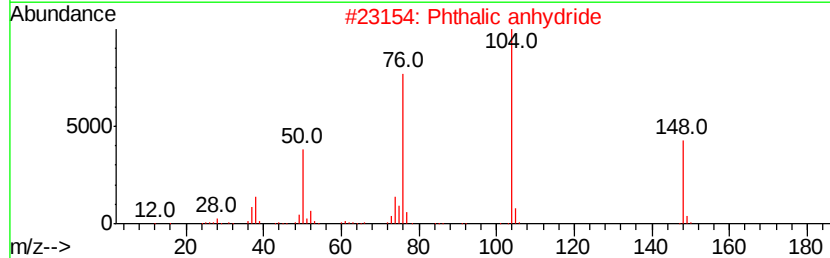
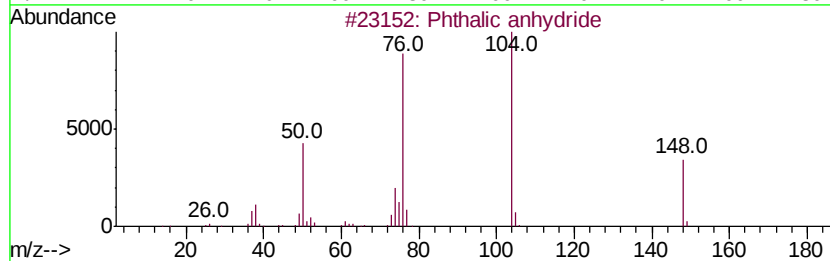
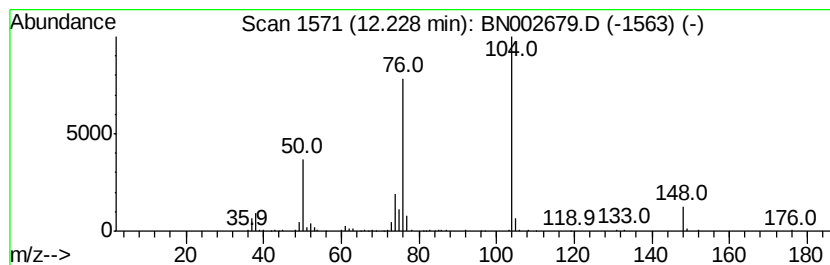
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Phthalic anhydride Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	4.14 ng/ul	397277	Naphthalene-d8	10.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic anhydride	148	C8H4O3	000085-44-9	95
2		Phthalic anhydride	148	C8H4O3	000085-44-9	95
3		Phthalic anhydride	148	C8H4O3	000085-44-9	90
4		[2,2'-Bi-1H-indene]-1,1'-dione, ...	358	C18H14O8	005950-69-6	90
5		Ninhydrin	178	C9H6O4	000485-47-2	90



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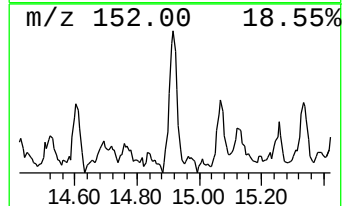
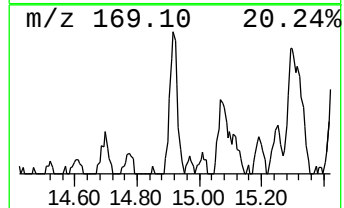
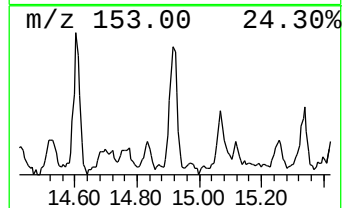
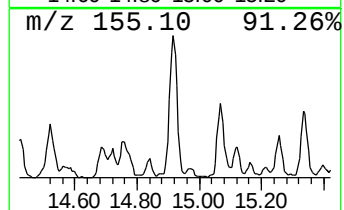
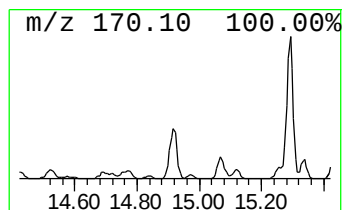
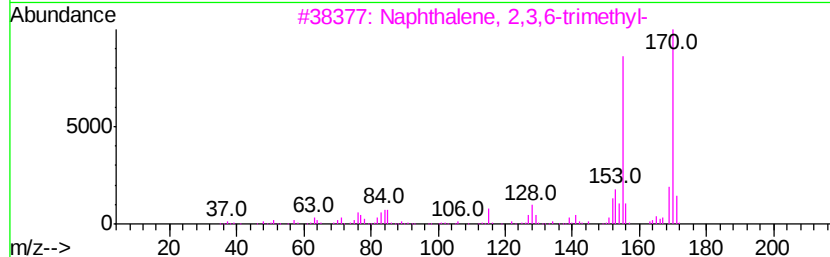
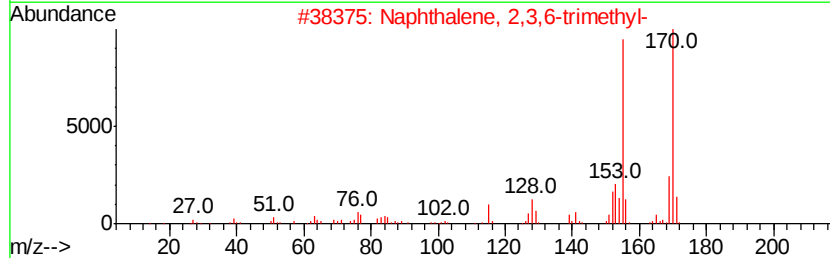
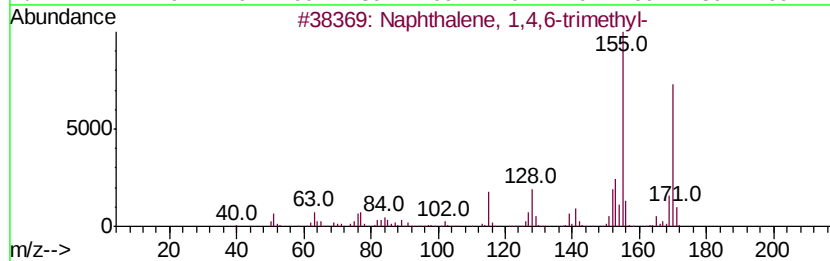
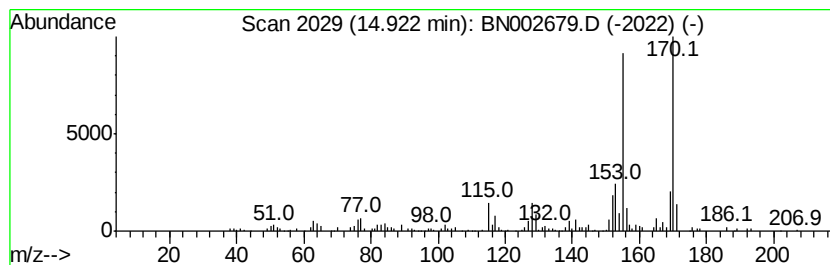
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Naphthalene, 1,4,6-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	2.66 ng/ul	332677	Acenaphthene-d10	14.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
4		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
5		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96



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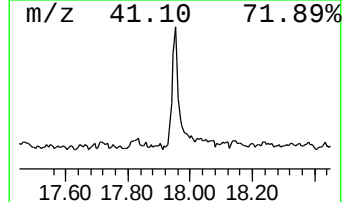
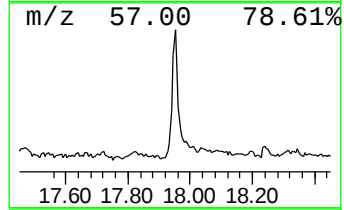
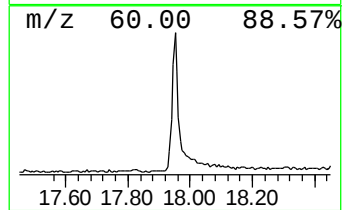
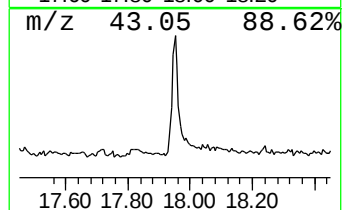
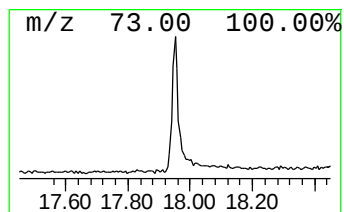
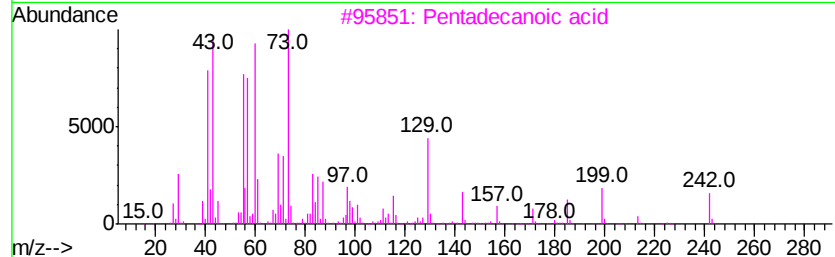
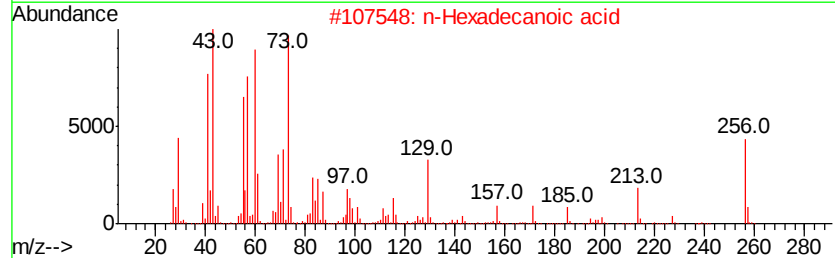
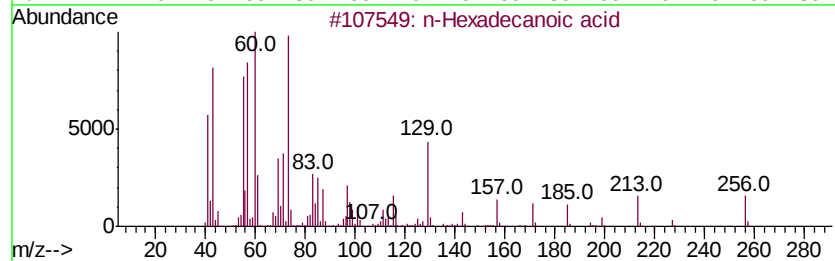
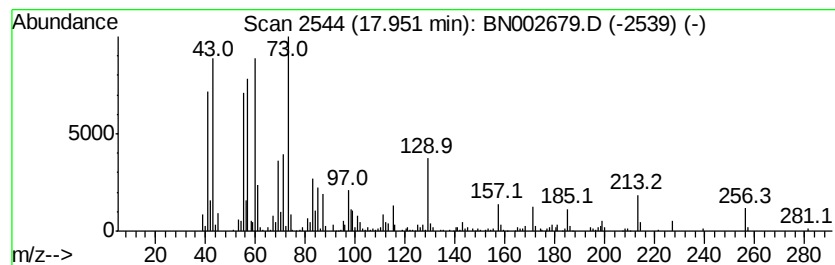
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.95	3.77 ng/ul	449243	Phenanthrene-d10		17.05	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	81
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
Data File : BN002679.D
Acq On : 13 Sep 2018 17:17
Operator : JU/SJ
Sample : J4864-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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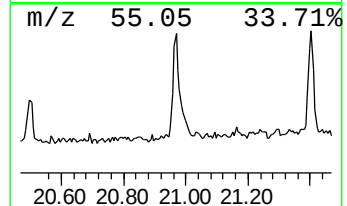
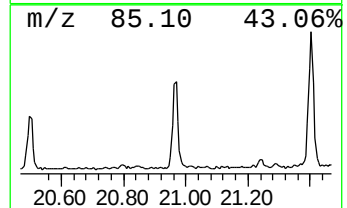
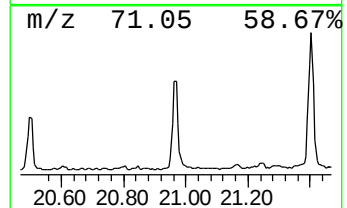
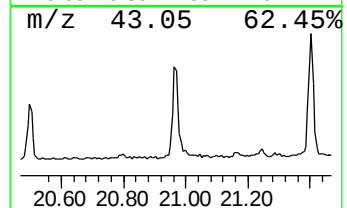
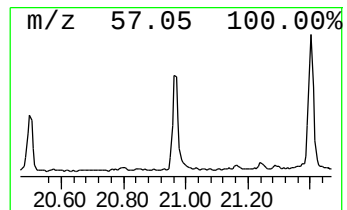
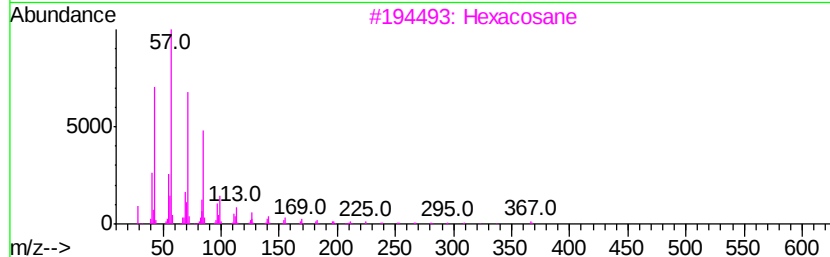
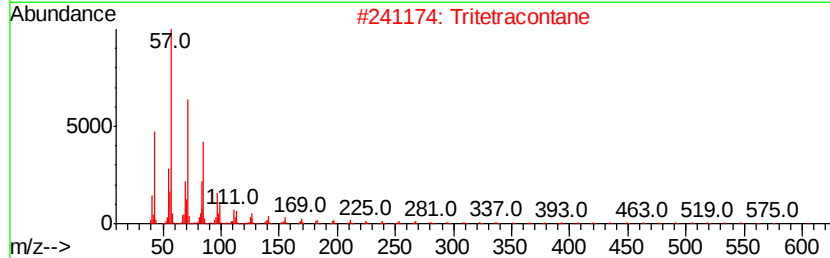
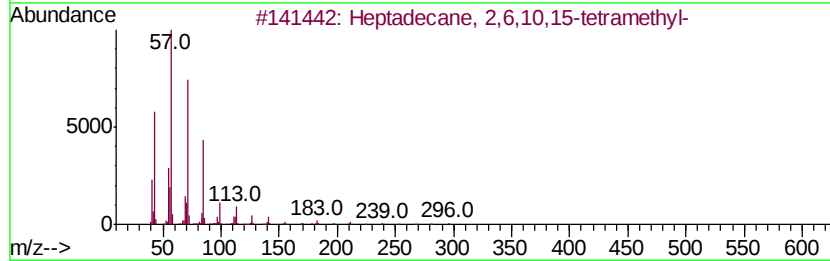
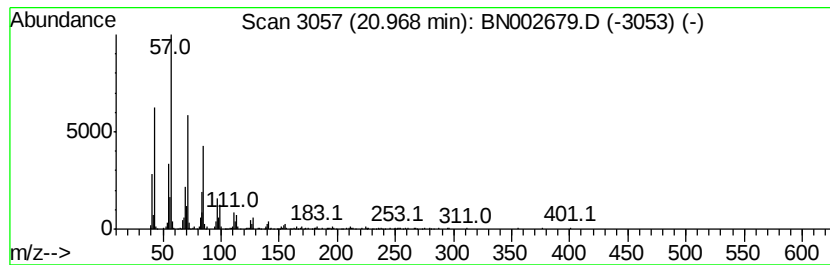
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	4.15 ng/ul	438817	Chrysene-d12	21.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	96
2	Tritetracontane	605	C43H88	007098-21-7	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Tetratetracontane	619	C44H90	007098-22-8	90
5	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
Data File : BN002679.D
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Instrument :
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ClientSampleId :
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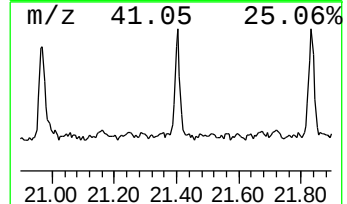
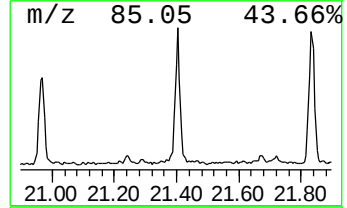
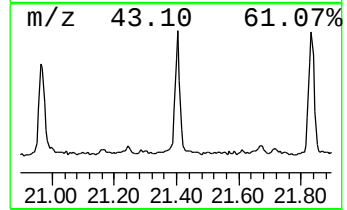
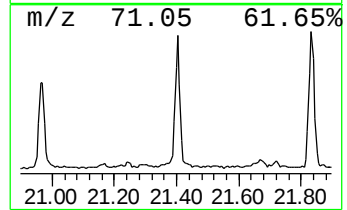
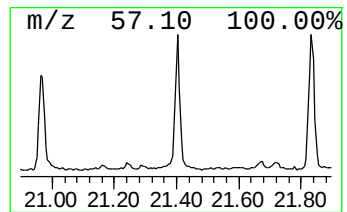
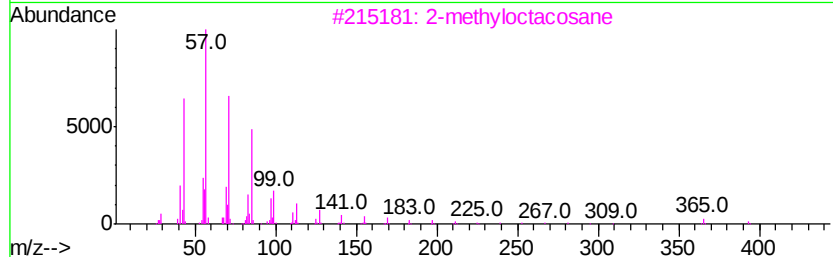
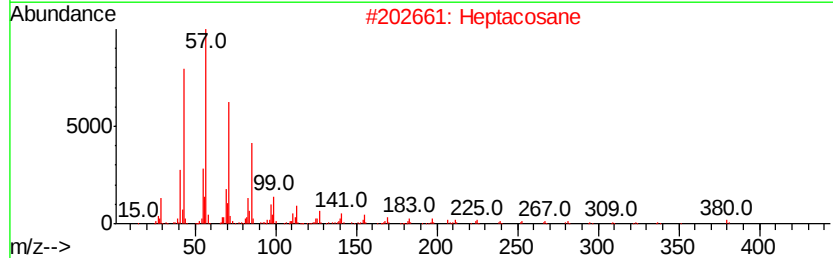
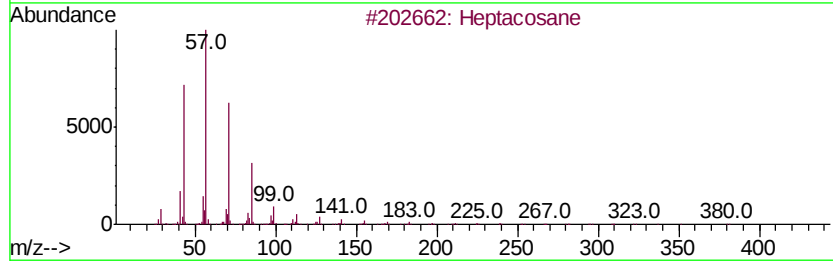
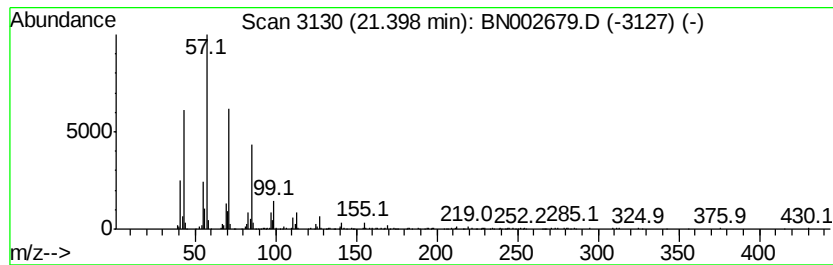
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.40	3.78 ng/ul	400201	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
Data File : BN002679.D
Acq On : 13 Sep 2018 17:17
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Sample : J4864-05
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Instrument :
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ClientSampleId :
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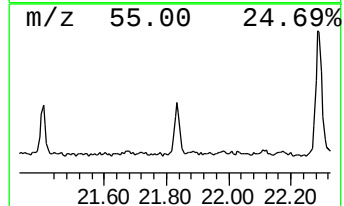
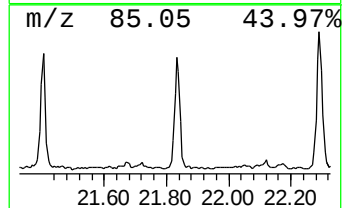
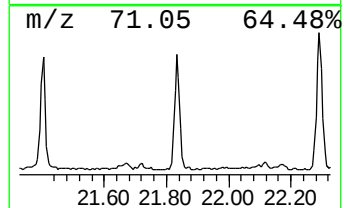
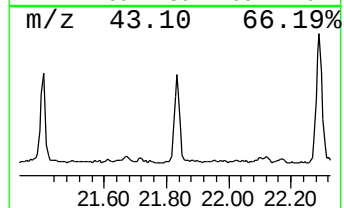
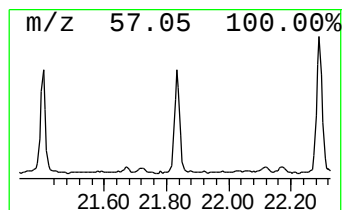
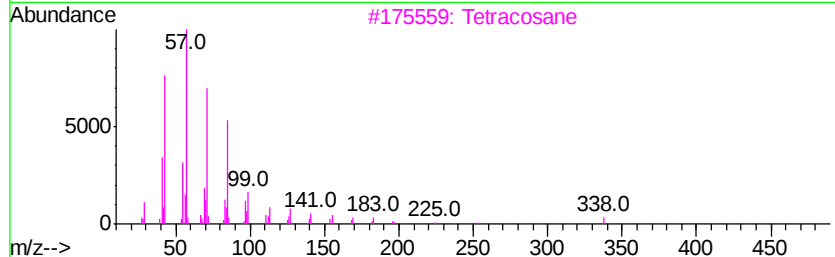
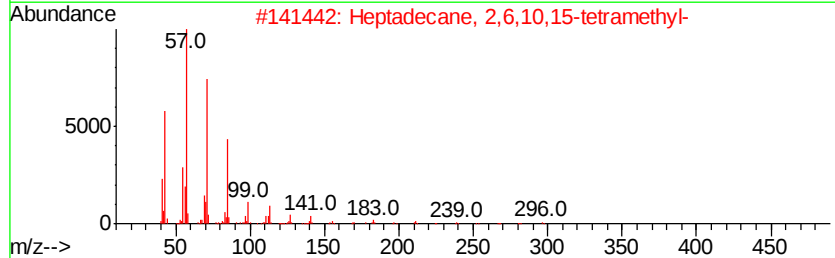
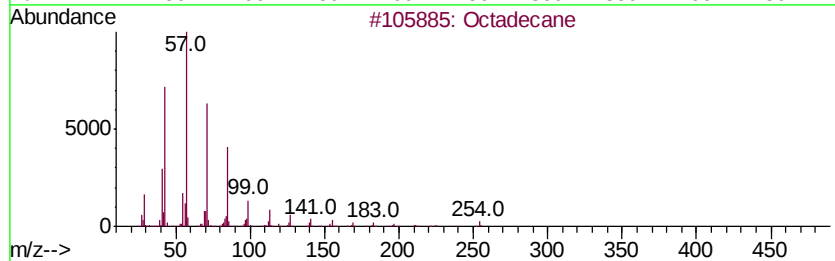
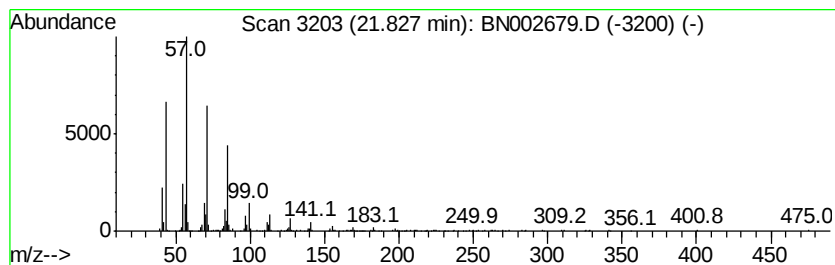
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	4.19 ng/ul	443000	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Tetracosane	338	C24H50	000646-31-1	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
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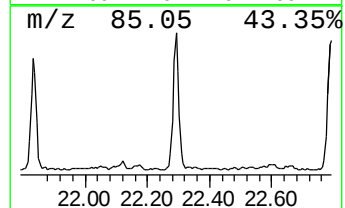
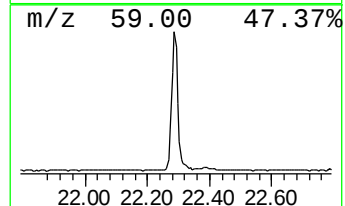
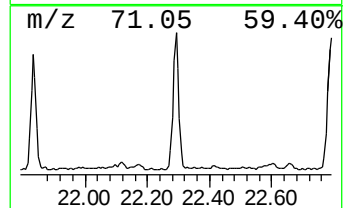
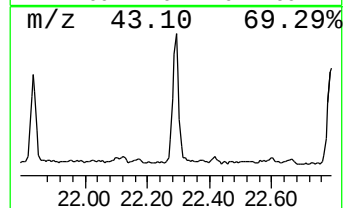
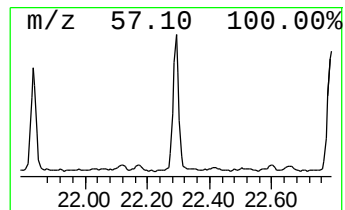
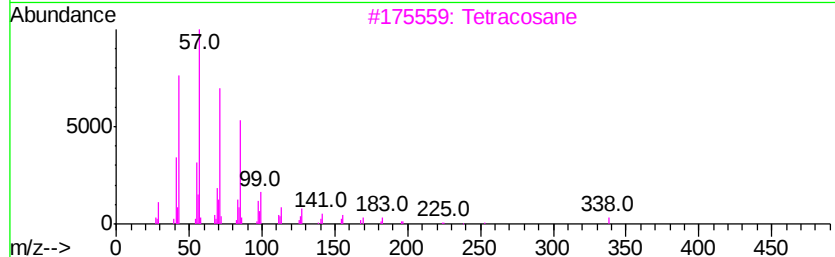
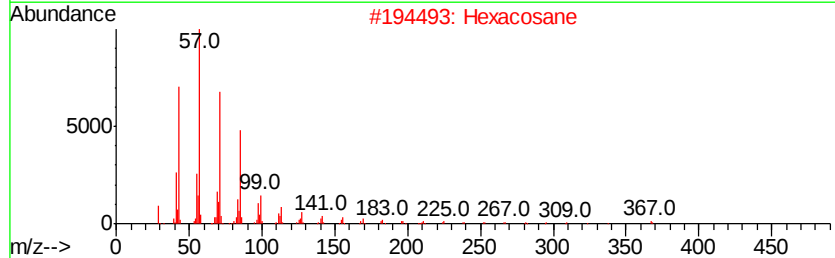
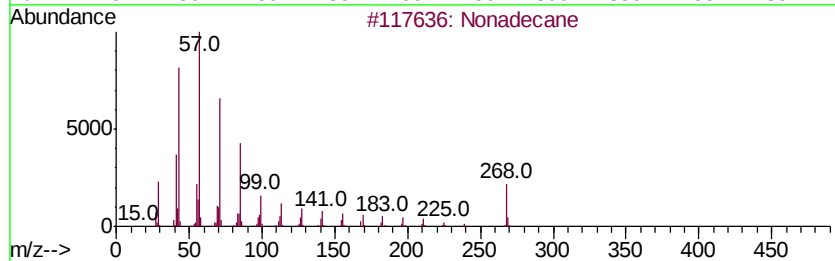
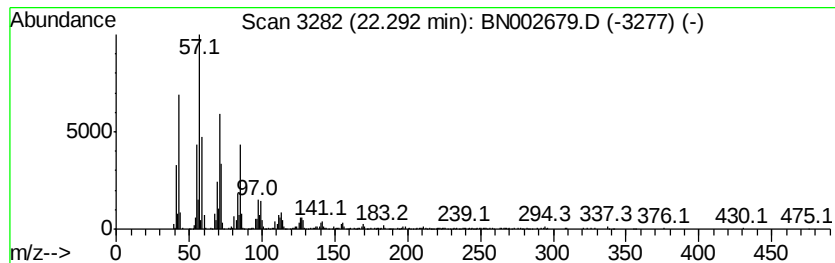
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.29	10.49 ng/ul	1108960	Chrysene-d12	21.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Hexacosane	366	C26H54	000630-01-3	64
3		Tetracosane	338	C24H50	000646-31-1	64
4		Tritetracontane	605	C43H88	007098-21-7	64
5		Octacosane	394	C28H58	000630-02-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
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Misc :
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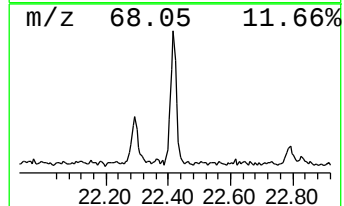
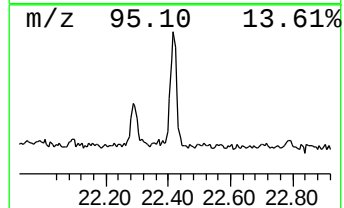
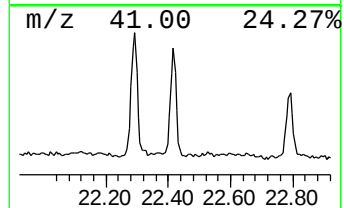
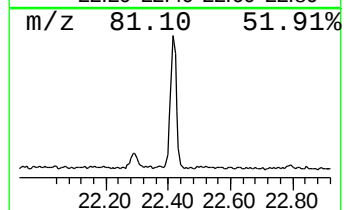
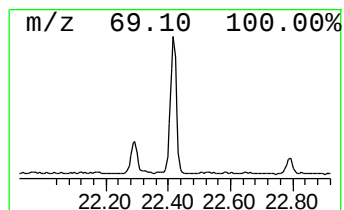
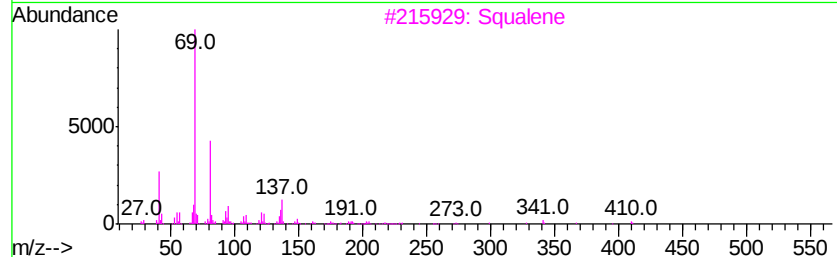
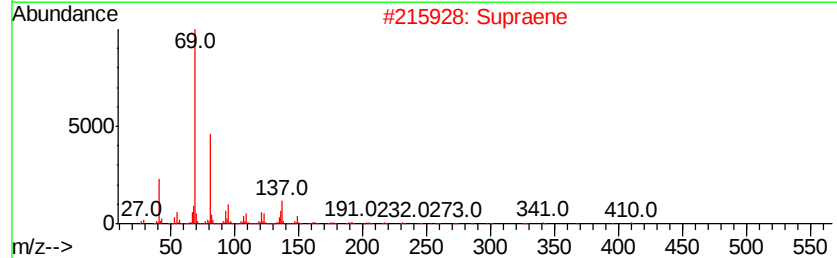
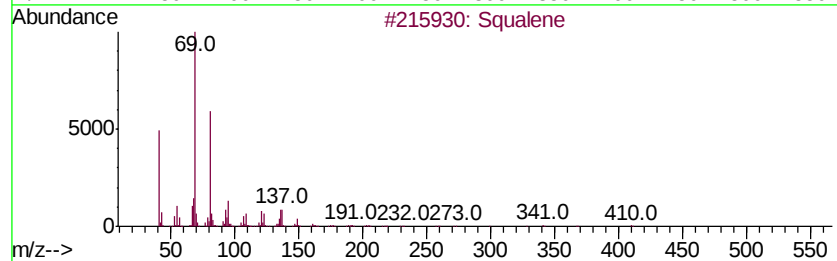
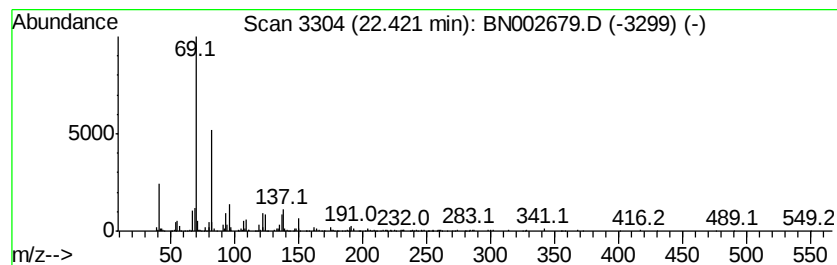
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Squalene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.42	5.12 ng/ul	552468	Perylene-d12	23.47
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	000111-02-4	87
2	Supraene	410 C30H50	007683-64-9	87
3	Squalene	410 C30H50	000111-02-4	83
4	7,11-Dimethyldodeca-2,6,10-trien...	208 C14H24O	125529-10-4	72
5	4,9,13,17-Tetramethyl-4,8,12,16-...	316 C22H36O	056882-09-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
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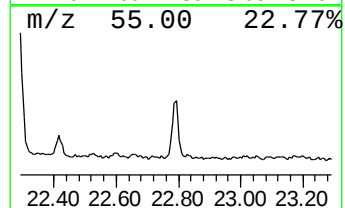
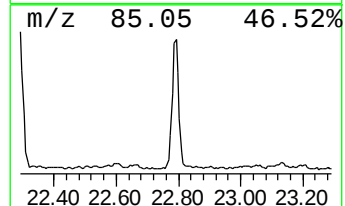
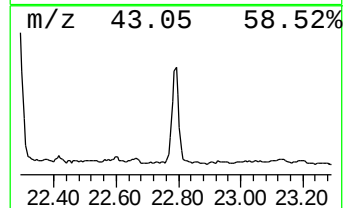
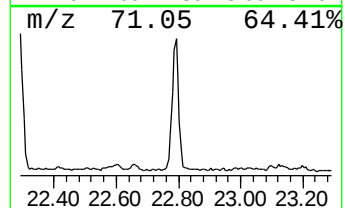
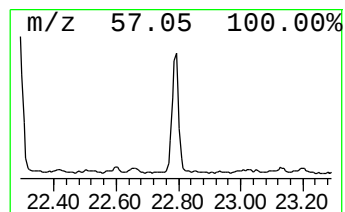
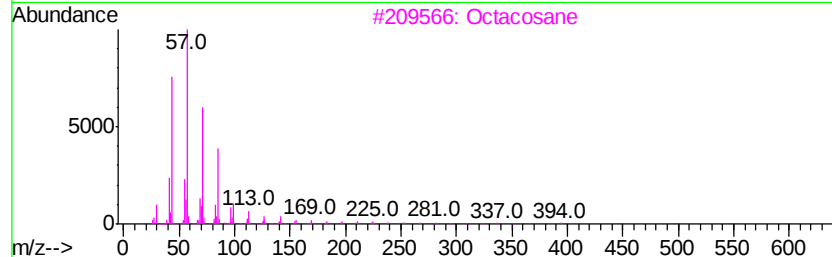
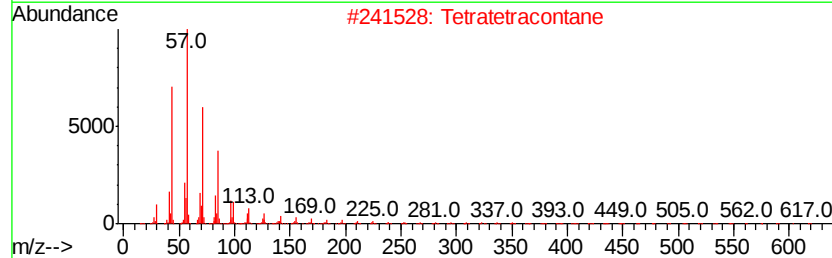
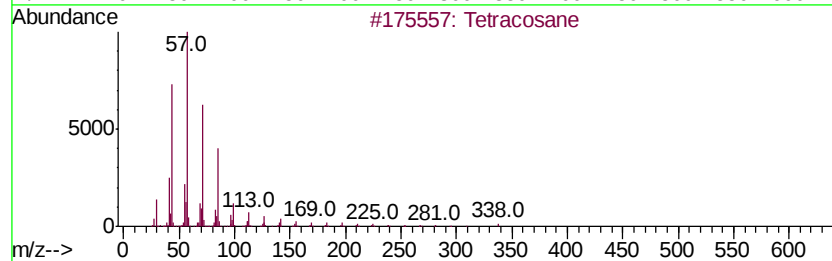
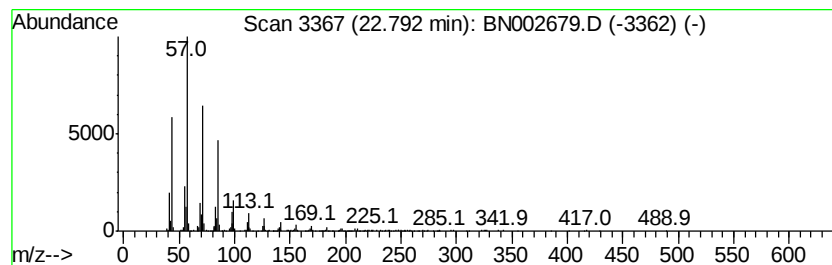
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	6.17 ng/ul	664804	Perylene-d12	23.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Octacosane	394	C28H58	000630-02-4	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
Data File : BN002679.D
Acq On : 13 Sep 2018 17:17
Operator : JU/SJ
Sample : J4864-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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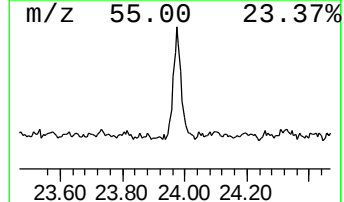
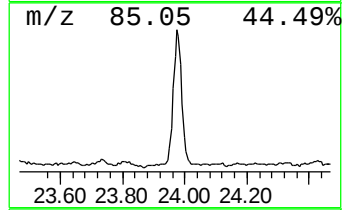
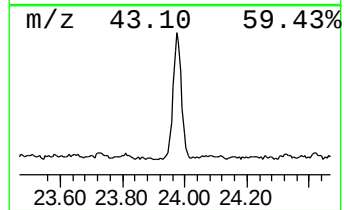
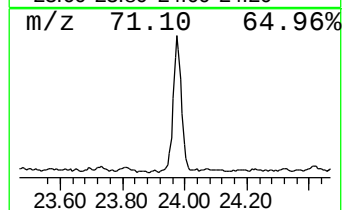
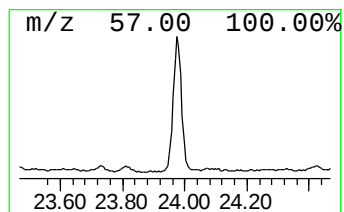
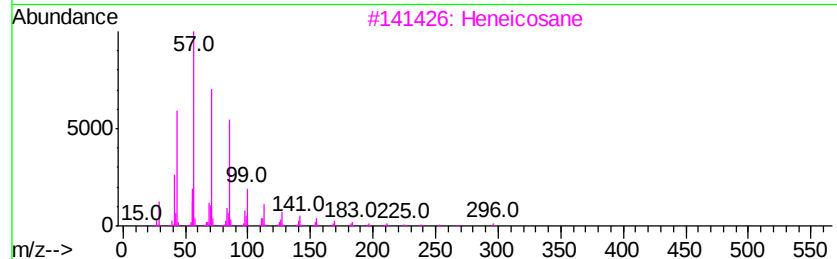
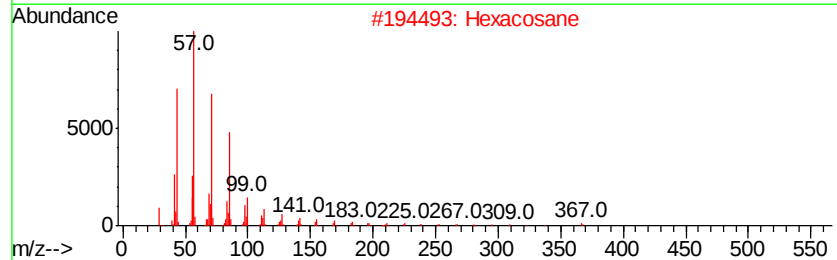
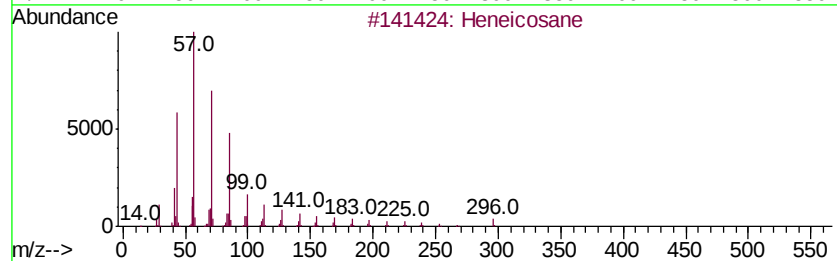
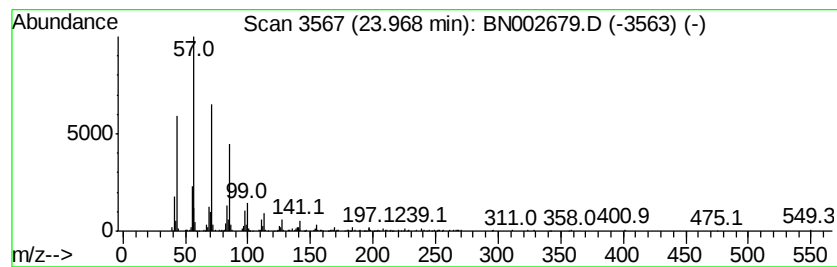
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.97	6.63 ng/ul	715100	Perylene-d12	23.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	93
2			Hexacosane	366	C26H54	000630-01-3	90
3			Heneicosane	296	C21H44	000629-94-7	90
4			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN091318\
Data File : BN002679.D
Acq On : 13 Sep 2018 17:17
Operator : JU/SJ
Sample : J4864-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0D11

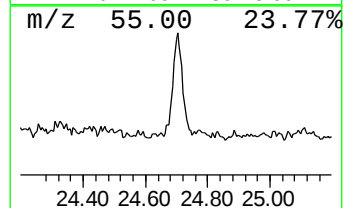
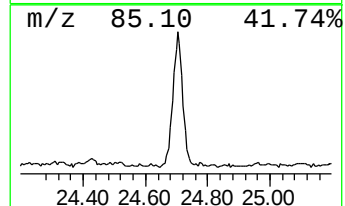
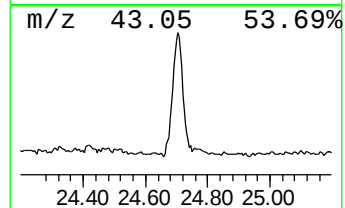
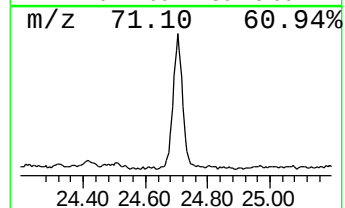
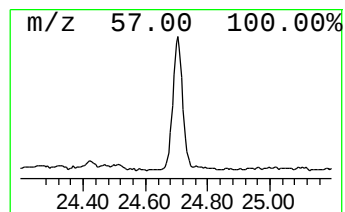
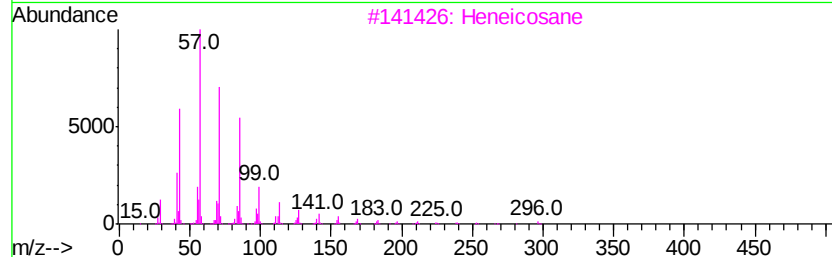
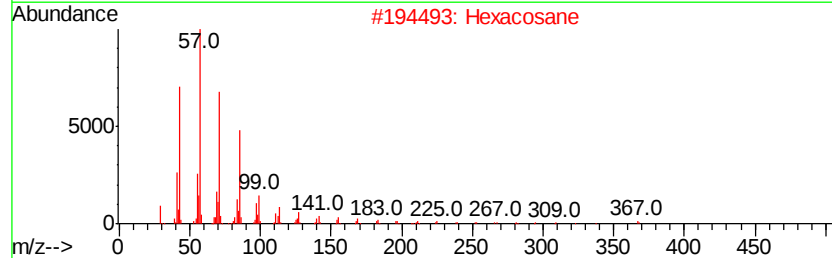
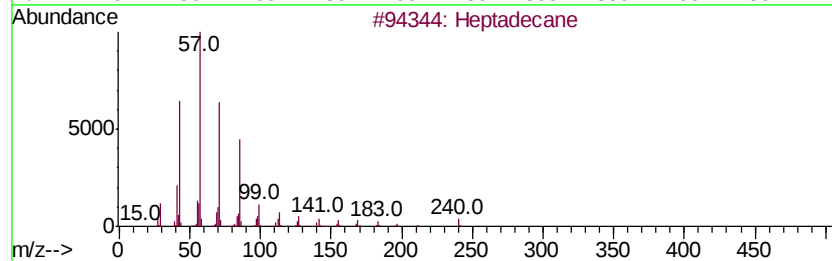
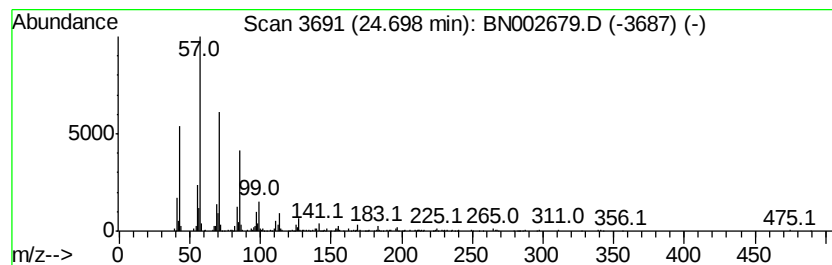
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.70	5.02 ng/ul	541089	Perylene-d12	23.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	94
2			Hexacosane	366	C26H54	000630-01-3	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN091318\
Data File : BN002679.D
Acq On : 13 Sep 2018 17:17
Operator : JU/SJ
Sample : J4864-05
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0D11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN081318MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzyl alcohol	7.91	4.5	ng/ul	271077	1	7.66	1196530	20.0
Phthalic anhydride	12.23	4.1	ng/ul	397277	2	10.43	1918200	20.0
Naphthalene, 1,4,...	14.92	2.7	ng/ul	332677	3	14.29	2504100	20.0
n-Hexadecanoic acid	17.95	3.8	ng/ul	449243	4	17.05	2382300	20.0
(DEL) Alkane: Str...	20.97	4.2	ng/ul	438817	5	21.24	2115050	20.0
(DEL) Alkane: Str...	21.40	3.8	ng/ul	400201	5	21.24	2115050	20.0
(DEL) Alkane: Str...	21.83	4.2	ng/ul	443000	5	21.24	2115050	20.0
(DEL) Alkane: Str...	22.29	10.5	ng/ul	1108960	5	21.24	2115050	20.0
Squalene	22.42	5.1	ng/ul	552468	6	23.47	2156530	20.0
(DEL) Alkane: Str...	22.79	6.2	ng/ul	664804	6	23.47	2156530	20.0
(DEL) Alkane: Str...	23.97	6.6	ng/ul	715100	6	23.47	2156530	20.0
(DEL) Alkane: Str...	24.70	5.0	ng/ul	541089	6	23.47	2156530	20.0