

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
 Data File : BP000784.D
 Acq On : 15 Oct 2019 22:36
 Operator : JU
 Sample : K5178-09
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPF2

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_P\METHODS\SOM-EPA-BP100919MA.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	2.358	42	46	56	rBV	32281	47354	1.45%	0.098%
2	3.505	234	241	250	rBV	22172	38581	1.19%	0.080%
3	3.716	273	277	280	rVV	12378	19283	0.59%	0.040%
4	3.752	280	283	290	rVB	15973	22603	0.69%	0.047%
5	4.916	478	481	485	rVB	13149	13103	0.40%	0.027%
6	6.305	711	717	723	rBV	11595	17779	0.55%	0.037%
7	6.387	727	731	736	rBV	1034953	1001712	30.77%	2.082%
8	6.434	736	739	750	rVV	851928	801822	24.63%	1.667%
9	6.522	750	754	761	rVV	1122197	1031405	31.68%	2.144%
10	6.593	761	766	776	rVB3	15658	27273	0.84%	0.057%
11	6.752	789	793	796	rBV	1361320	1235653	37.95%	2.569%
12	6.816	801	804	811	rVB	10675	11737	0.36%	0.024%
13	7.134	854	858	871	rBV	1260018	1066738	32.77%	2.218%
14	7.304	884	887	891	rVV	1055896	895302	27.50%	1.861%
15	7.522	921	924	928	rVB	25883	21453	0.66%	0.045%
16	7.634	939	943	955	rBV	949605	746933	22.94%	1.553%
17	7.881	982	985	994	rBV	1295430	1233968	37.90%	2.565%
18	7.940	994	995	1007	rVV3	20638	34629	1.06%	0.072%
19	8.034	1007	1011	1014	rVV	2223349	1715035	52.68%	3.565%
20	8.093	1018	1021	1036	rVB	1039522	871750	26.78%	1.812%
21	8.646	1110	1115	1117	rBV2	16021	15674	0.48%	0.033%
22	8.716	1124	1127	1131	rBV	24932	21205	0.65%	0.044%
23	9.234	1211	1215	1218	rVV2	18360	19740	0.61%	0.041%
24	9.451	1248	1252	1255	rVV2	13210	13996	0.43%	0.029%
25	9.487	1255	1258	1260	rVV	2080705	1696419	52.11%	3.527%
26	9.510	1260	1262	1267	rVV	1193858	877216	26.94%	1.824%
27	9.640	1281	1284	1288	rBV	2842793	2255139	69.27%	4.688%
28	9.798	1307	1311	1314	rBV	2536062	2178381	66.91%	4.529%
29	9.910	1327	1330	1339	rBV	405657	478539	14.70%	0.995%
30	9.975	1339	1341	1344	rVV	60924	58705	1.80%	0.122%
31	10.004	1344	1346	1353	rVB2	142834	121147	3.72%	0.252%
32	10.210	1378	1381	1384	rVV2	11543	13060	0.40%	0.027%
33	10.316	1396	1399	1403	rBV	3025763	2694971	82.78%	5.603%
34	10.345	1403	1404	1408	rVB	49683	28838	0.89%	0.060%

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35	10.393	1408	1412	1418	rBV	534308	456154	14.01%	0.948%
36	10.445	1418	1421	1429	rVV	48462	56002	1.72%	0.116%
37	10.628	1448	1452	1460	rVV6	5530	11329	0.35%	0.024%
38	10.716	1464	1467	1470	rVV2	12995	18038	0.55%	0.037%
39	10.745	1470	1472	1475	rVB	17324	13985	0.43%	0.029%
40	10.834	1484	1487	1491	rVB	13697	11872	0.36%	0.025%
41	11.034	1518	1521	1524	rBV	16272	13004	0.40%	0.027%
42	11.069	1524	1527	1530	rVB	53474	41997	1.29%	0.087%
43	11.104	1530	1533	1537	rVB2	9665	11294	0.35%	0.023%
44	11.169	1537	1544	1546	rBV2	21552	28425	0.87%	0.059%
45	11.257	1555	1559	1561	rBV3	20469	20014	0.61%	0.042%
46	11.292	1561	1565	1568	rBV	2724155	2392856	73.50%	4.975%
47	11.316	1568	1569	1571	rVB	104320	59682	1.83%	0.124%
48	11.351	1571	1575	1578	rBV	3400923	2743569	84.27%	5.704%
49	11.528	1600	1605	1613	rBV2	24635	40524	1.24%	0.084%
50	11.698	1630	1634	1639	rVB3	15975	17924	0.55%	0.037%
51	11.840	1654	1658	1660	rBV2	97295	106020	3.26%	0.220%
52	11.863	1660	1662	1667	rVB	42159	44940	1.38%	0.093%
53	11.916	1668	1671	1673	rBV	17881	16896	0.52%	0.035%
54	12.010	1685	1687	1690	rBV2	19218	17745	0.55%	0.037%
55	12.104	1700	1703	1706	rBV2	21591	21030	0.65%	0.044%
56	12.228	1721	1724	1726	rBV	14546	12086	0.37%	0.025%
57	12.357	1744	1746	1749	rBV	12206	12618	0.39%	0.026%
58	12.404	1751	1754	1758	rVB3	39389	40967	1.26%	0.085%
59	12.498	1768	1770	1771	rBV	36107	28410	0.87%	0.059%
60	12.522	1771	1774	1777	rVB	178596	152405	4.68%	0.317%
61	12.634	1790	1793	1795	rBV3	10545	13849	0.43%	0.029%
62	12.657	1795	1797	1804	rVB2	38099	42135	1.29%	0.088%
63	12.734	1806	1810	1816	rVV	3395319	3255598	100.00%	6.768%
64	12.781	1816	1818	1820	rVV	45102	38469	1.18%	0.080%
65	12.804	1820	1822	1827	rVB2	34922	30790	0.95%	0.064%
66	12.898	1835	1838	1845	rVB2	27044	36877	1.13%	0.077%
67	13.086	1868	1870	1873	rVB	28482	22296	0.68%	0.046%
68	13.145	1877	1880	1884	rVV2	65144	67259	2.07%	0.140%
69	13.192	1885	1888	1895	rVB2	19544	29991	0.92%	0.062%
70	13.287	1901	1904	1906	rBV	20540	18176	0.56%	0.038%
71	13.316	1907	1909	1913	rVB5	18069	16890	0.52%	0.035%

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72	13.381	1917	1920	1924	rVV	53719	54168	1.66%	0.113%
73	13.492	1936	1939	1941	rBV	79290	62350	1.92%	0.130%
74	13.739	1979	1981	1983	rBV	30736	24681	0.76%	0.051%
75	13.775	1983	1987	1991	rBV3	89620	112969	3.47%	0.235%
76	13.822	1991	1995	1999	rBV2	319338	450113	13.83%	0.936%
77	13.863	1999	2002	2008	rBV	1971224	1804269	55.42%	3.751%
78	13.922	2009	2012	2015	rBV	225223	182102	5.59%	0.379%
79	13.969	2015	2020	2023	rBV2	2477208	3038719	93.34%	6.317%
80	14.051	2031	2034	2035	rBV	165717	151916	4.67%	0.316%
81	14.069	2035	2037	2040	rVV	211316	236595	7.27%	0.492%
82	14.110	2040	2044	2046	rVV2	813474	974120	29.92%	2.025%
83	14.128	2046	2047	2048	rVV	591947	357838	10.99%	0.744%
84	14.145	2048	2050	2057	rVV	541542	837190	25.72%	1.740%
85	14.198	2057	2059	2063	rVV	186191	190910	5.86%	0.397%
86	14.251	2066	2068	2073	rVV	328736	459061	14.10%	0.954%
87	14.292	2073	2075	2076	rVV	407131	347101	10.66%	0.722%
88	14.310	2076	2078	2085	rVV	646318	1076831	33.08%	2.239%
89	14.363	2085	2087	2091	rVB2	158815	136961	4.21%	0.285%
90	14.457	2097	2103	2107	rBV2	307406	432972	13.30%	0.900%
91	14.575	2121	2123	2133	rVB	65941	97017	2.98%	0.202%
92	14.763	2152	2155	2159	rBV	234221	224251	6.89%	0.466%
93	15.016	2195	2198	2206	rBV	86149	150887	4.63%	0.314%
94	15.104	2209	2213	2216	rBV	270602	307470	9.44%	0.639%
95	15.245	2234	2237	2244	rVB	58403	84853	2.61%	0.176%
96	15.357	2251	2256	2266	rVB	1942507	2391220	73.45%	4.971%
97	15.445	2266	2271	2275	rVV	1615325	1922275	59.05%	3.996%
98	15.486	2275	2278	2290	rVB2	204410	288648	8.87%	0.600%
99	15.922	2348	2352	2358	rBV	171216	254872	7.83%	0.530%
100	16.433	2435	2439	2452	rVB2	85155	160741	4.94%	0.334%

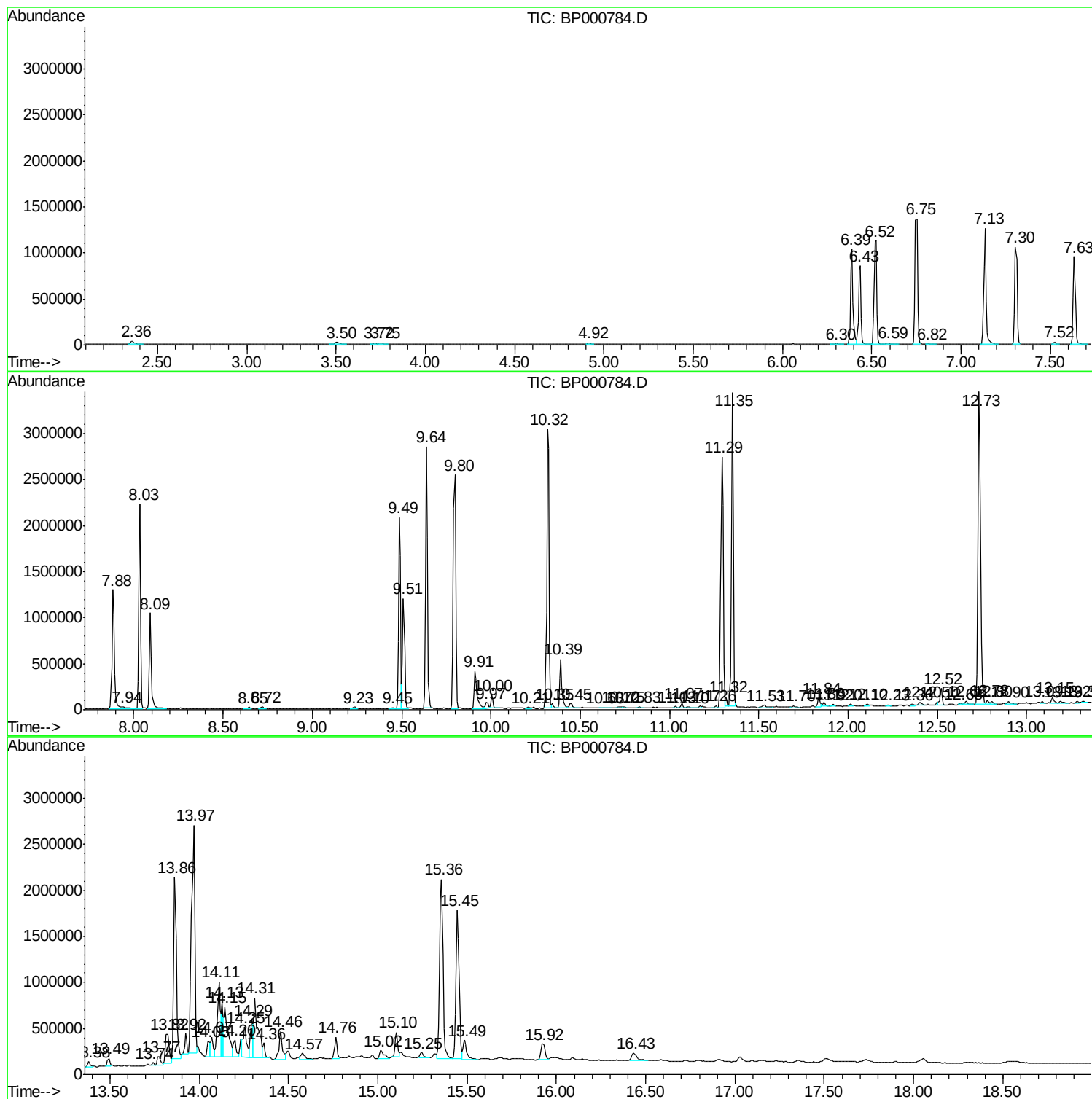
Sum of corrected areas: 48102359

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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



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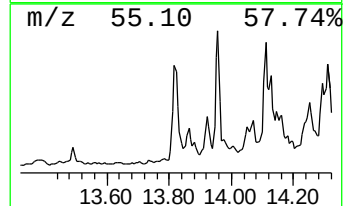
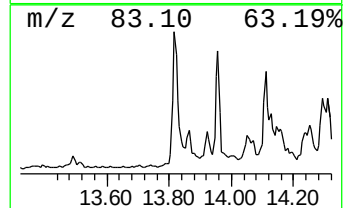
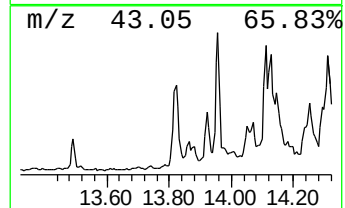
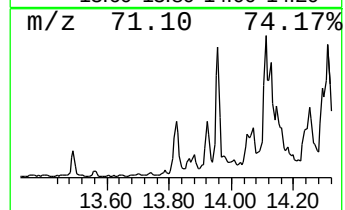
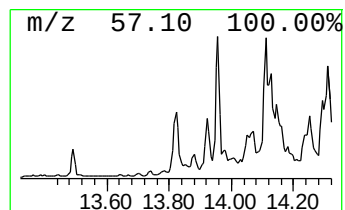
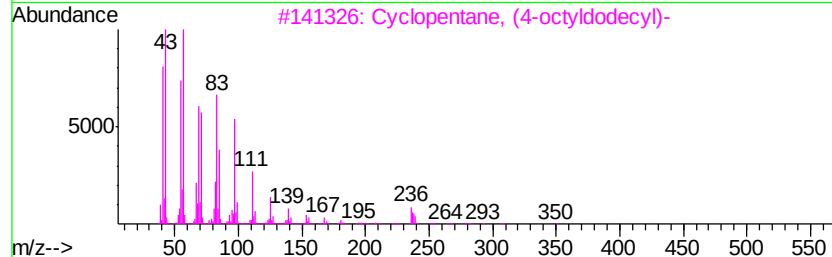
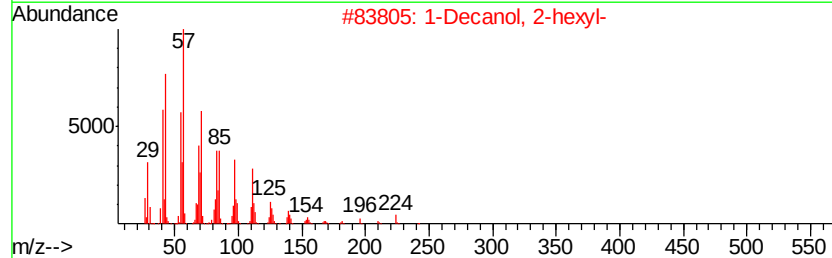
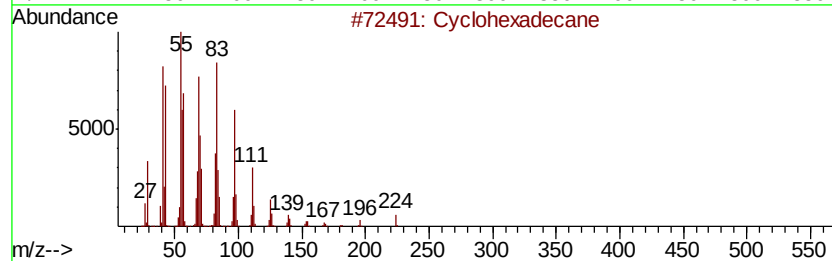
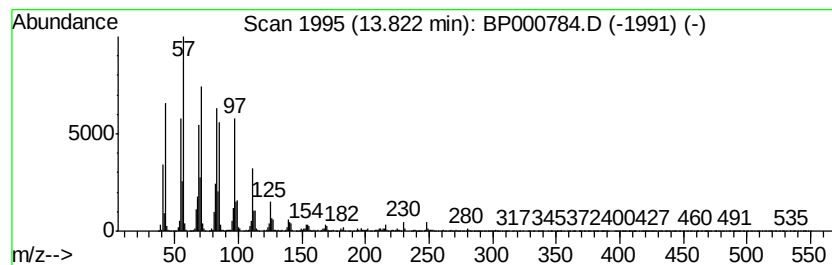
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic13.82 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	2.96 ng/ul	450113	Chrysene-d12	13.97
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 90
2		1-Decanol, 2-hexyl-	242 C16H34O	002425-77-6 86
3		Cyclopentane, (4-octylidodecyl)-	350 C25H50	005638-09-5 74
4		Heptafluorobutanoic acid, hepta...	452 C21H35F7O2	1000282-97-3 70
5		Pentafluoropropionic acid, hepta...	402 C20H35F5O2	1000283-04-2 70



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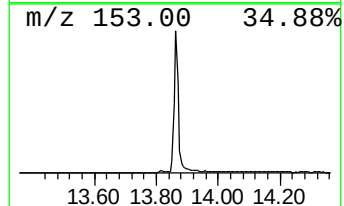
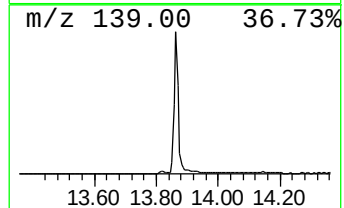
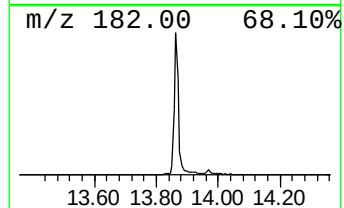
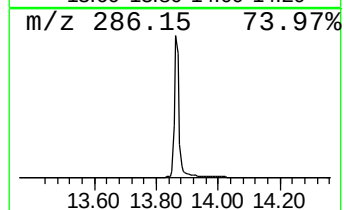
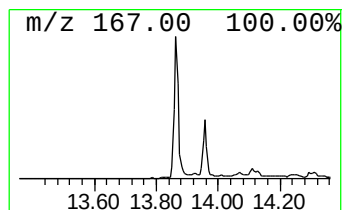
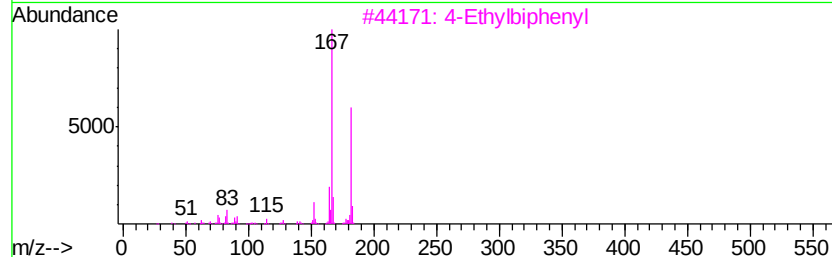
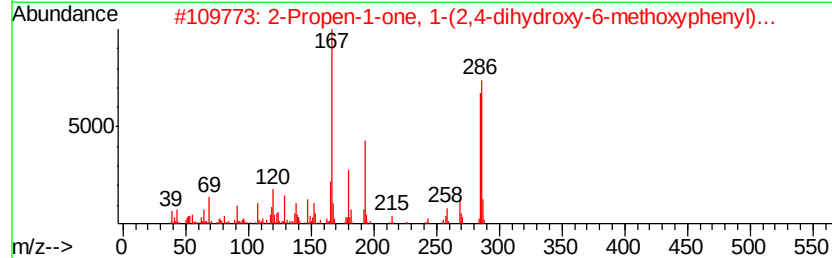
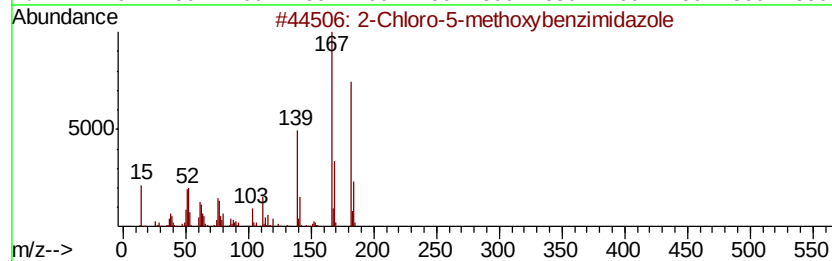
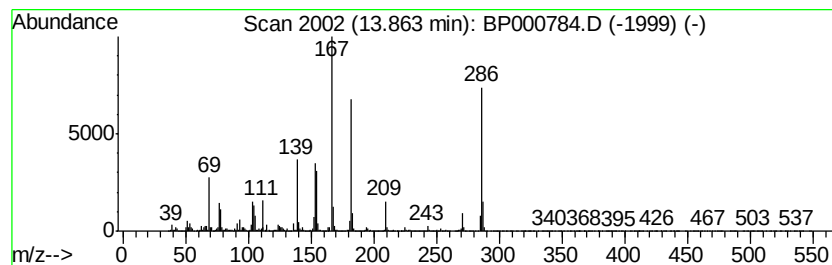
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	11.88 ng/ul	1804270	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Chloro-5-methoxybenzimidazole	182	C8H7ClN2O	015965-54-5	43
2		2-Propen-1-one, 1-(2,4-dihydroxy...	286	C16H14O5	069707-17-1	43
3		4-Ethylbiphenyl	182	C14H14	005707-44-8	38
4		1,7-Dinitrophenazine 5-oxide	286	C12H6N4O5	108272-86-2	38
5		Benzene, 1-methyl-4-(phenylmethyl)-	182	C14H14	000620-83-7	35



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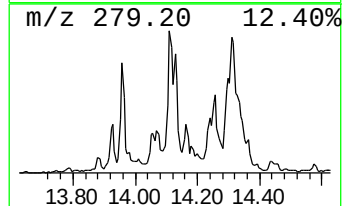
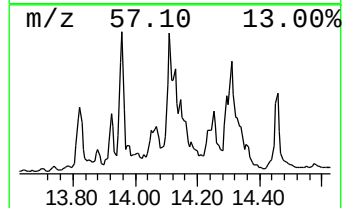
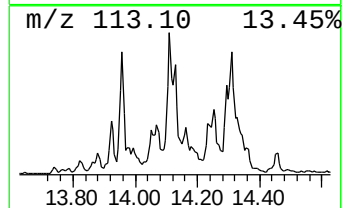
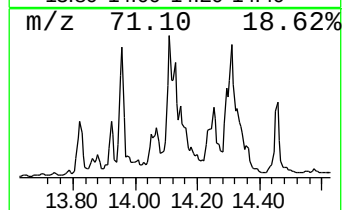
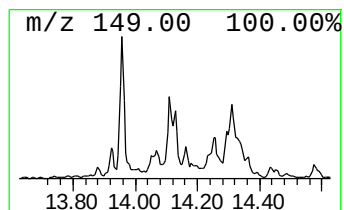
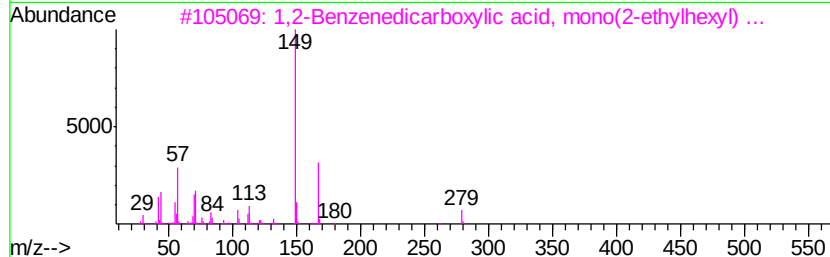
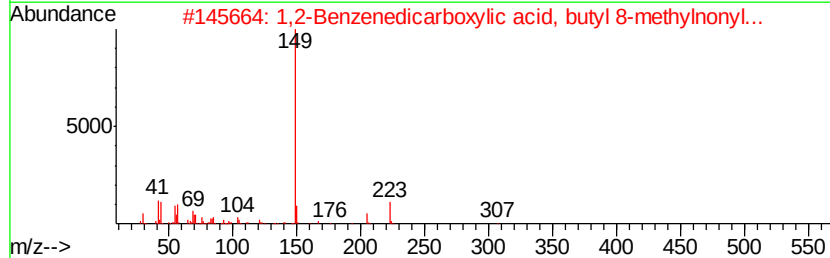
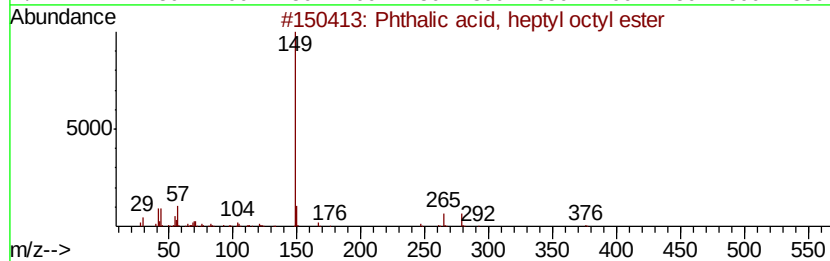
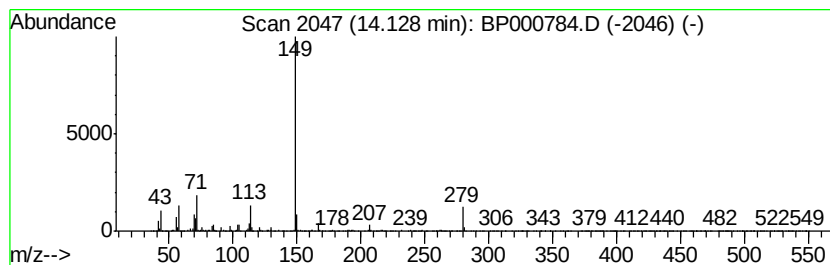
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Peak Number 4 Phthalic acid, heptyl octyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	2.36 ng/ul	357838	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, heptyl octyl ester	376	C23H36O4	088216-58-4	59
2		1,2-Benzenedicarboxylic acid, bu...	362	C22H34O4	000089-18-9	53
3		1,2-Benzenedicarboxylic acid, mo...	278	C16H22O4	004376-20-9	50
4		1,2-Benzenedicarboxylic acid, is...	418	C26H42O4	001330-96-7	50
5		1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	50



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Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
Operator : JU
Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

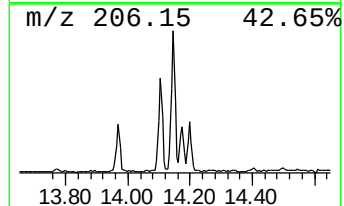
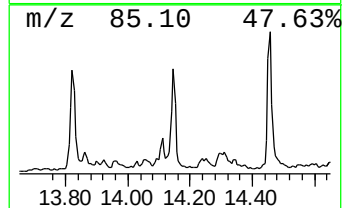
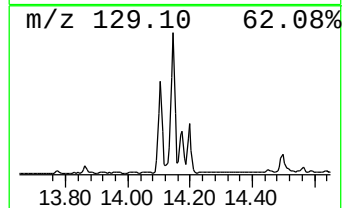
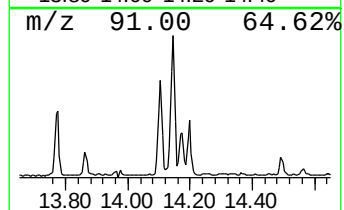
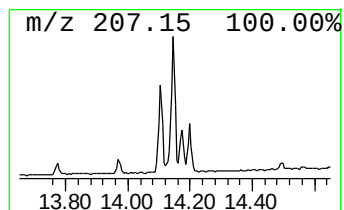
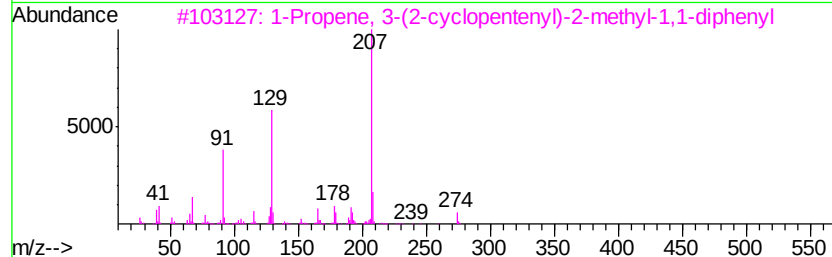
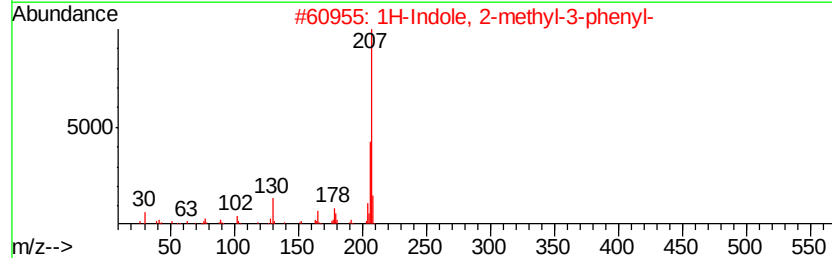
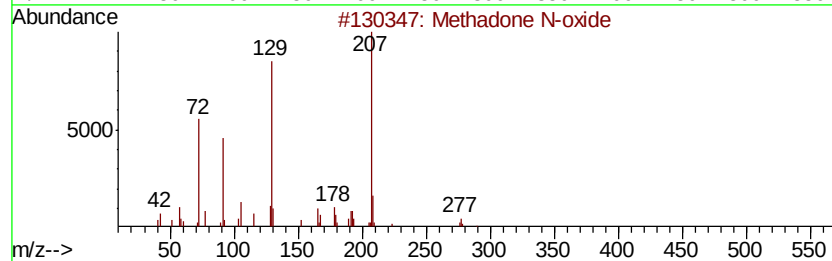
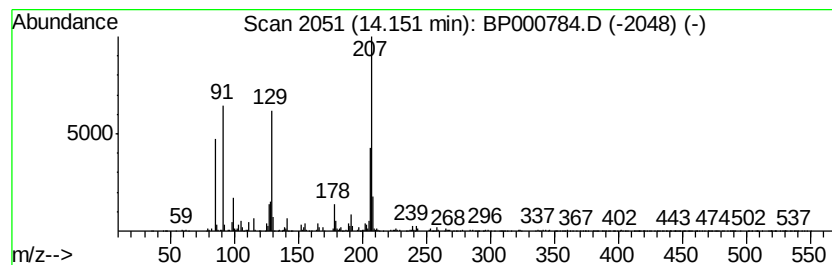
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Methadone N-oxide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.15	5.51 ng/ul	837190	Chrysene-d12	13.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methadone N-oxide	325	C21H27NO2	1000120-80-7	59
2		1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	41
3		1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	38
4		5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	38
5		1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
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Sample : K5178-09
Misc :
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Instrument :
BNA_P
ClientSampleId :
BEPF2

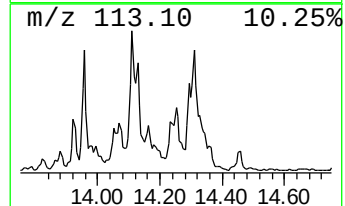
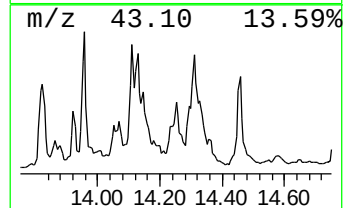
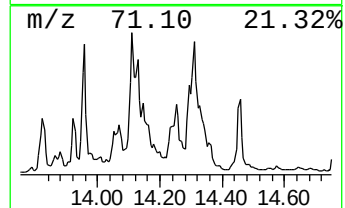
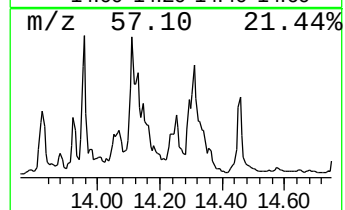
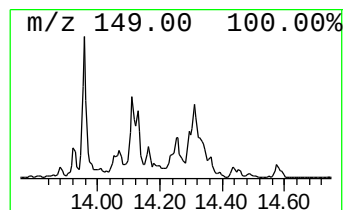
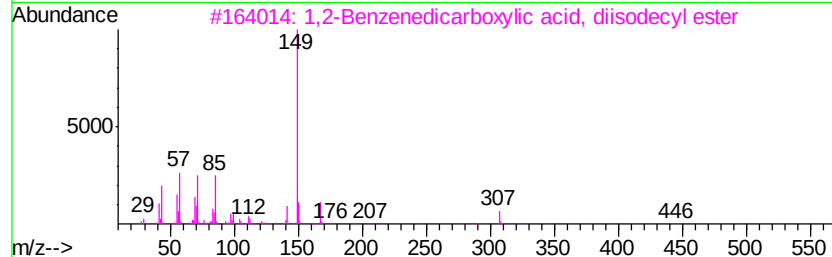
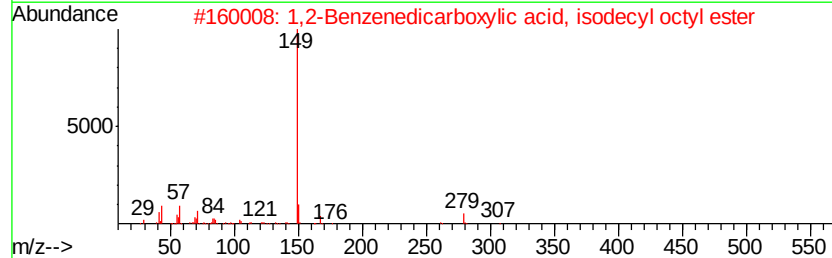
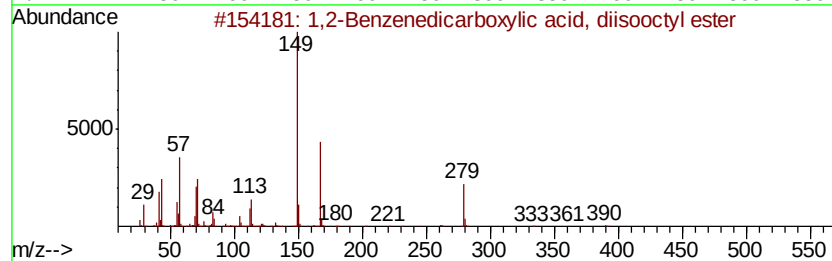
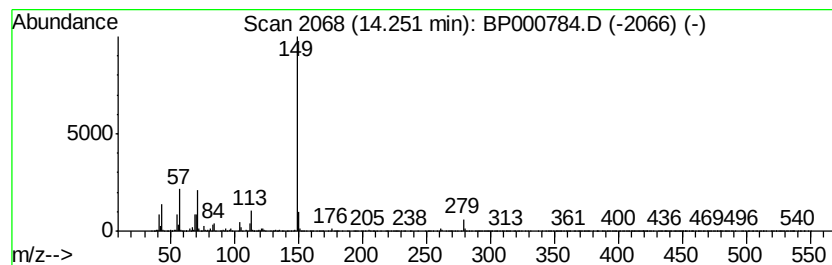
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,2-Benzenedicarboxylic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	3.02 ng/ul	459061	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	64
2		1,2-Benzenedicarboxylic acid, is...	418	C26H42O4	001330-96-7	64
3		1,2-Benzenedicarboxylic acid, di...	446	C28H46O4	026761-40-0	59
4		1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	53
5		1,2-Benzenedicarboxylic acid, bi...	334	C20H30O4	000146-50-9	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
Operator : JU
Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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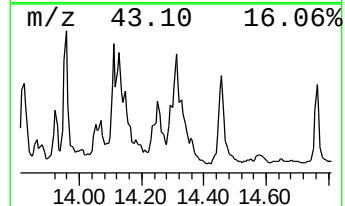
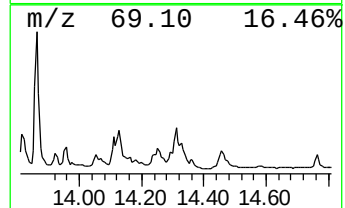
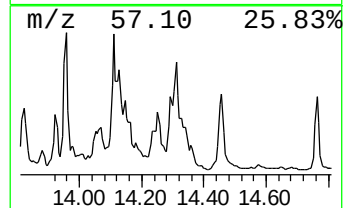
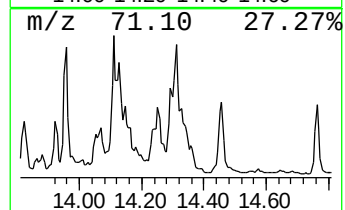
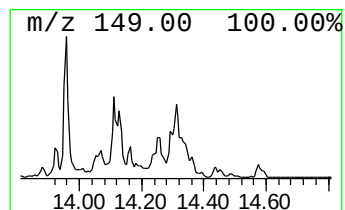
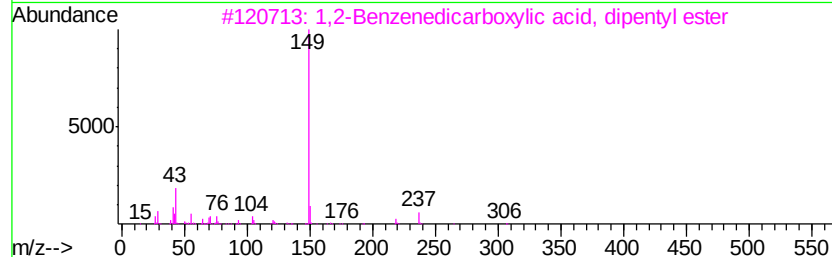
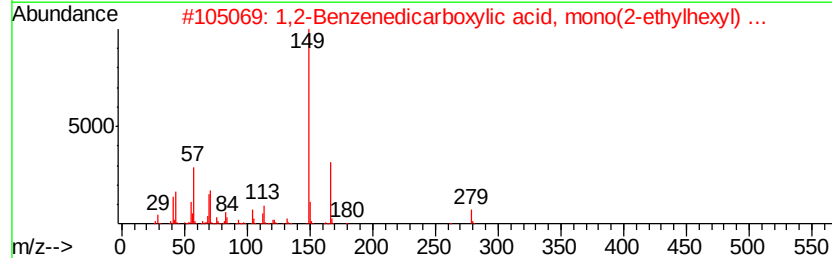
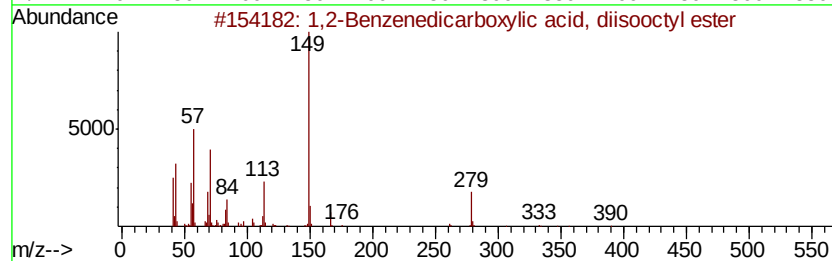
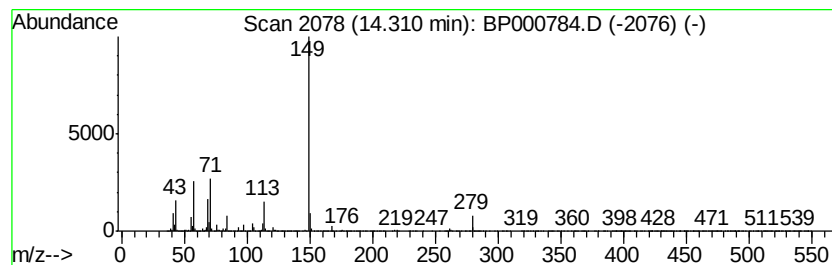
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,2-Benzenedicarboxylic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.31	7.09 ng/ul	1076830	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Benzenedicarboxylic acid, di...	390	C24H38O4	027554-26-3	83
2		1,2-Benzenedicarboxylic acid, mo...	278	C16H22O4	004376-20-9	64
3		1,2-Benzenedicarboxylic acid, di...	306	C18H26O4	000131-18-0	58
4		Phthalic acid, heptyl octyl ester	376	C23H36O4	088216-58-4	53
5		1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
Operator : JU
Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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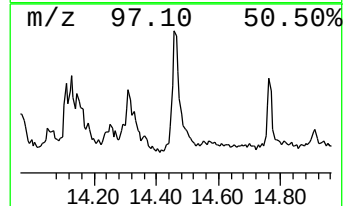
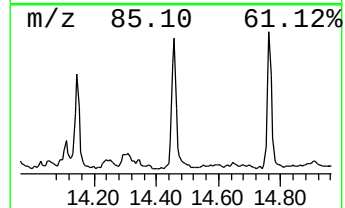
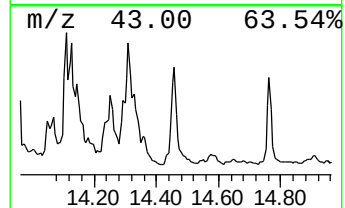
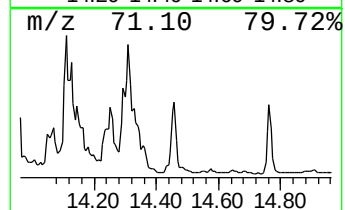
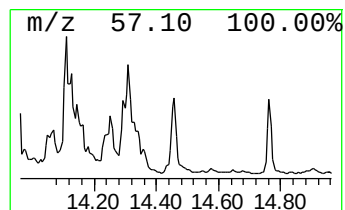
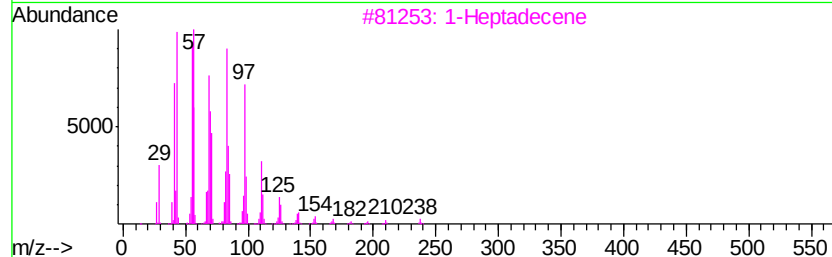
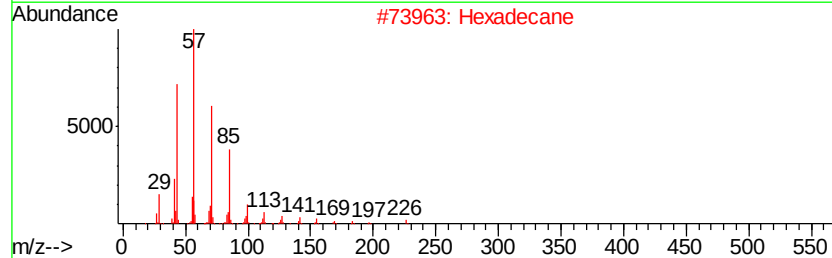
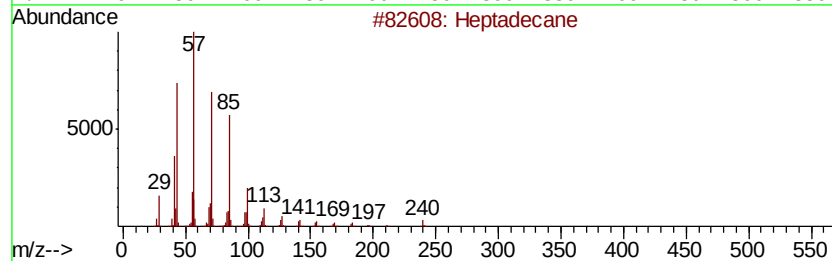
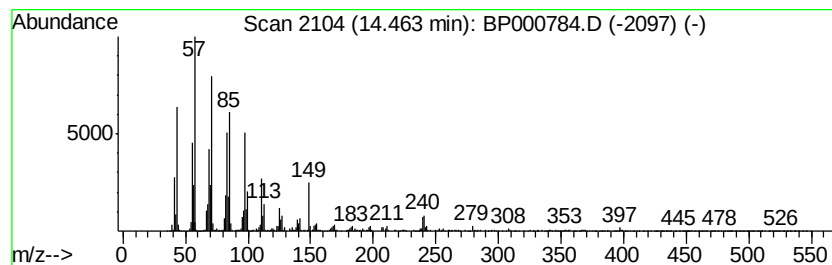
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	2.85 ng/ul	432972	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	96
2		Hexadecane	226	C16H34	000544-76-3	92
3		1-Heptadecene	238	C17H34	006765-39-5	92
4		Nonadecane	268	C19H40	000629-92-5	90
5		Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
Operator : JU
Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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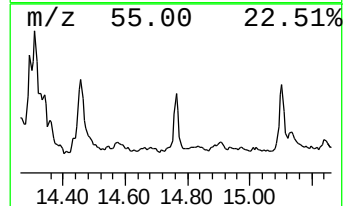
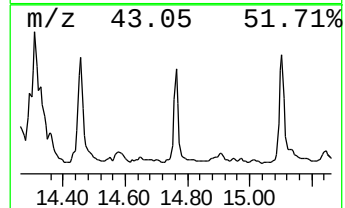
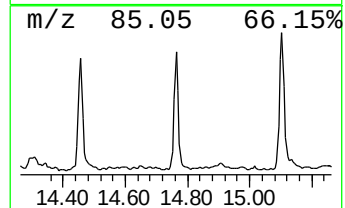
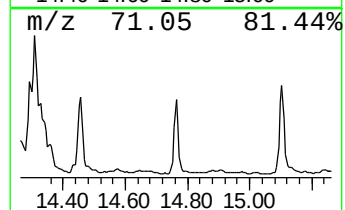
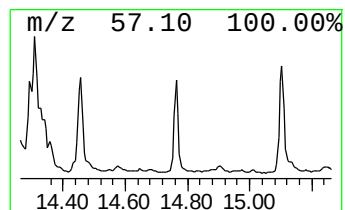
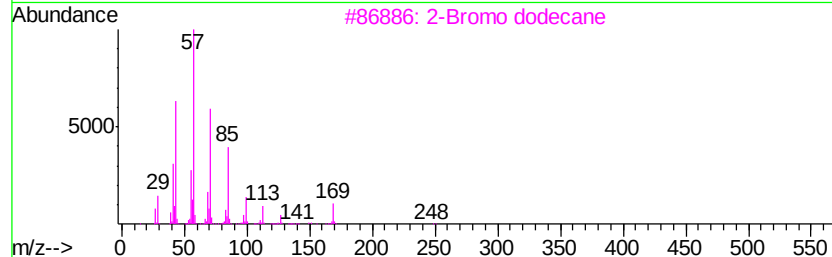
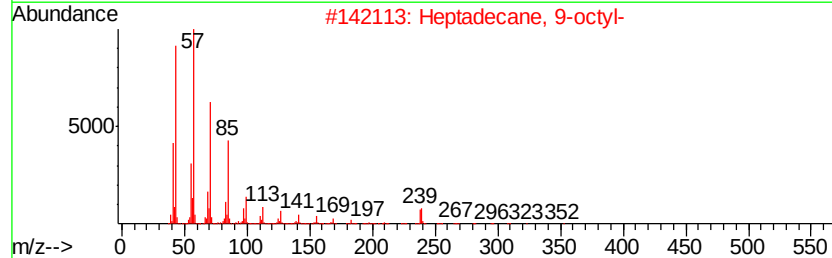
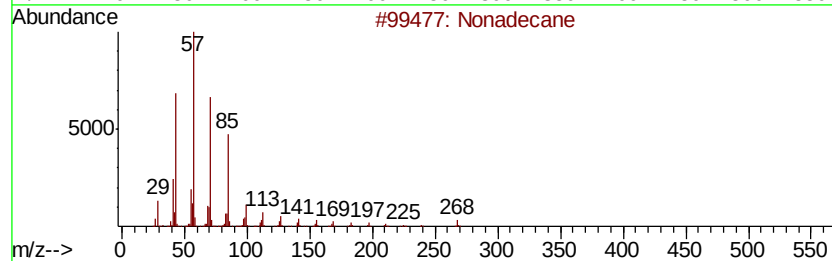
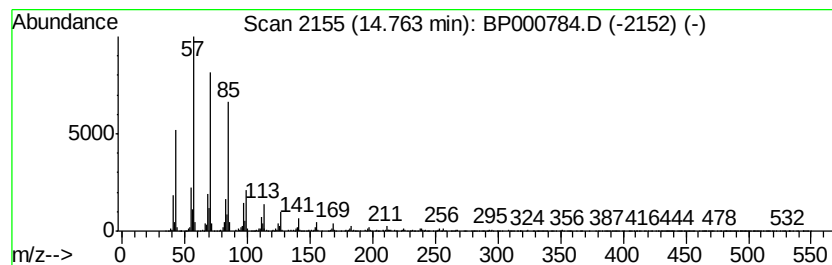
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	2.33 ng/ul	224251	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	95
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
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ClientSampleId :
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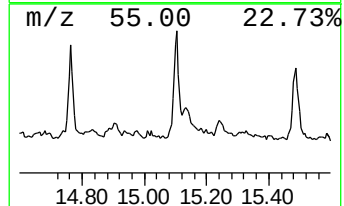
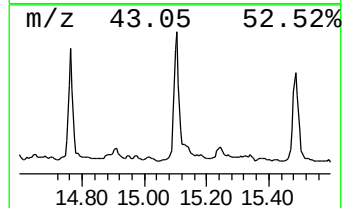
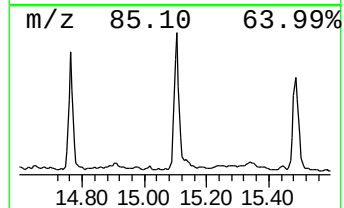
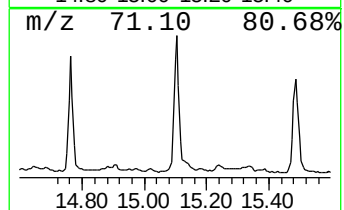
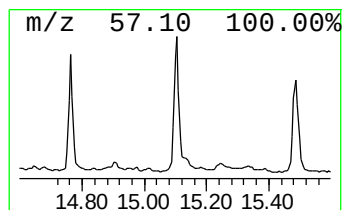
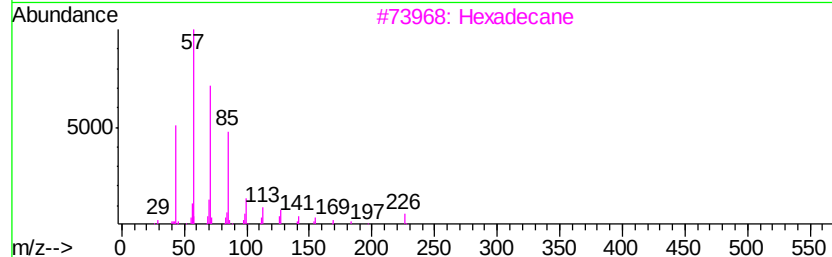
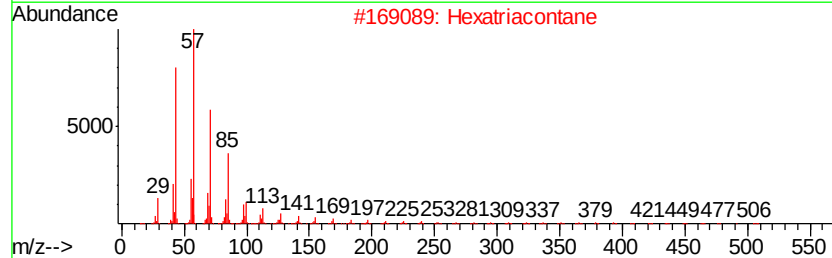
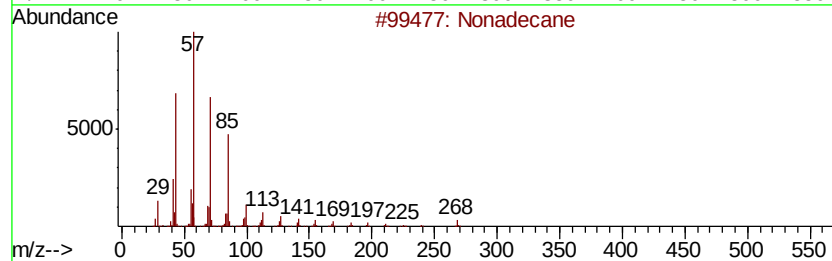
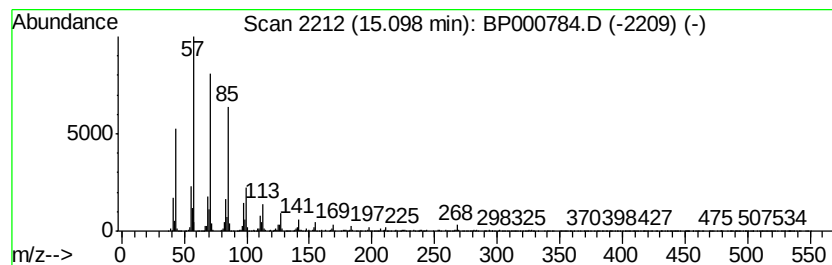
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	3.20 ng/ul	307470	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	94
2		Hexatriacontane	507	C36H74	000630-06-8	94
3		Hexadecane	226	C16H34	000544-76-3	93
4		Heptacosane	380	C27H56	000593-49-7	93
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
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Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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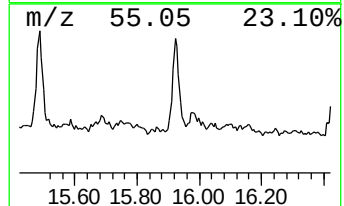
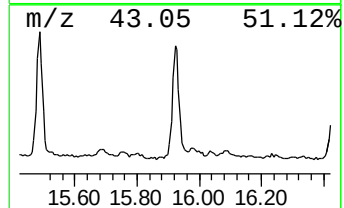
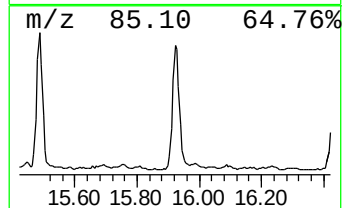
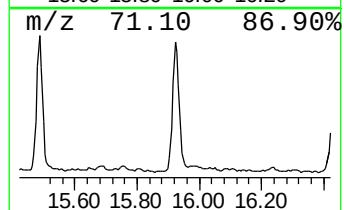
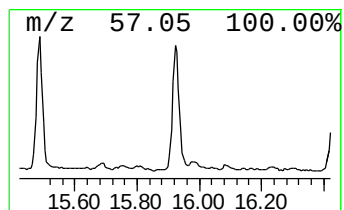
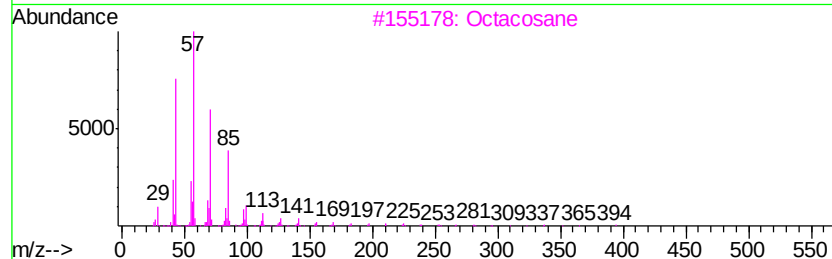
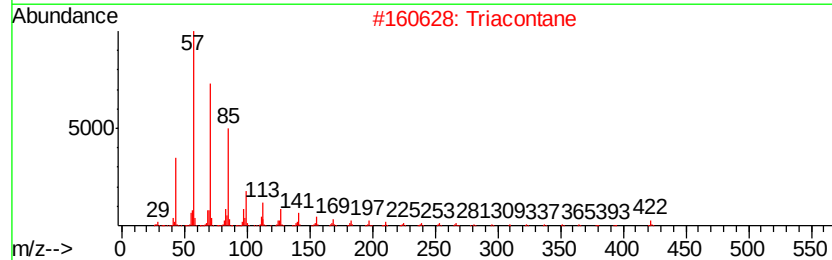
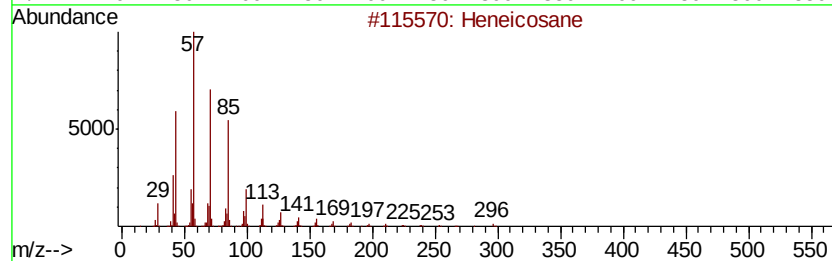
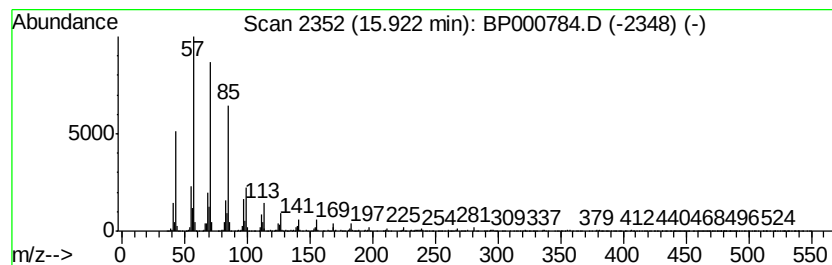
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	2.65 ng/ul	254872	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Triacontane	422	C30H62	000638-68-6	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Docosane, 9-butyl-	366	C26H54	055282-14-9	90
5		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101619\
Data File : BP000784.D
Acq On : 15 Oct 2019 22:36
Operator : JU
Sample : K5178-09
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BEPF2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.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
(DEL) Alkane: Cyc...	13.82	3.0	ng/ul	450113	5	13.97	3038720	20.0
unknown-01	13.86	11.9	ng/ul	1804270	5	13.97	3038720	20.0
Phthalic acid, he...	14.13	2.4	ng/ul	357838	5	13.97	3038720	20.0
Methadone N-oxide	14.15	5.5	ng/ul	837190	5	13.97	3038720	20.0
1,2-Benzenedicarb...	14.25	3.0	ng/ul	459061	5	13.97	3038720	20.0
1,2-Benzenedicarb...	14.31	7.1	ng/ul	1076830	5	13.97	3038720	20.0
(DEL) Alkane: Str...	14.46	2.9	ng/ul	432972	5	13.97	3038720	20.0
(DEL) Alkane: Str...	14.76	2.3	ng/ul	224251	6	15.45	1922280	20.0
(DEL) Alkane: Str...	15.10	3.2	ng/ul	307470	6	15.45	1922280	20.0
(DEL) Alkane: Str...	15.92	2.6	ng/ul	254872	6	15.45	1922280	20.0