LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000762.D
Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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_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	2.364	43	47	58	rBV	46998	65935	1.48%	0.099%
2	3.511	236	242	250	rBV	41603	69608	1.56%	0.105%
3	6.387	727	731	736	rBV2	1370192	1354139	30.37%	2.041%
4	6.434	736	739	750	rVV	1129264	1046037	23.46%	1.577%
5	6.522	750	754	757	rVV	1471437	1346684	30.20%	2.030%
6	6.746	789	792	796	rBV	1412859	1241219	27.84%	1.871%
7	7.134	854	858	873	rBV	1771312	1479283	33.18%	2.229%
8	7.305	884	887	891	rVV	1361393	1182965	26.53%	1.783%
9	7.634	939	943	946	rBV	1273086	961335	21.56%	1.449%
10	7.881	982	985	994	rBV	1719763	1619571	36.32%	2.441%
11	8.034	1007	1011	1014	rBV	2079348	1690332	37.91%	2.548%
12	8.093	1018	1021	1036	rVV	1398827	1160754	26.03%	1.749%
13	9.493	1255	1259	1261	rBV	2502282	2342070	52.53%	3.530%
14	9.510	1261	1262	1267	rVB	1193548	826323	18.53%	1.245%
15	9.640	1281	1284	1294	rVB	3333502	2939642	65.93%	4.430%
16	9.799	1307	1311	1314	rBV	2531583	2104127	47.19%	3.171%
17	9.828	1314	1316	1320	rVB	63769	49265	1.10%	0.074%
18	9.910	1326	1330	1339	rBV	654066	793854	17.80%	1.196%
19	9.981	1339	1342	1344	rVV	50697	52666	1.18%	0.079%
20	10.005	1344	1346	1351	rVV	110472	92359	2.07%	0.139%
21	10.210	1377	1381	1384	rBV2	31145	40844	0.92%	0.062%
22	10.322	1396	1400	1403	rBV	3626751	3372669	75.64%	5.083%
23	10.346	1403	1404	1409	rVB	115468	72319	1.62%	0.109%
24	10.393	1409	1412	1418	rBV	813257	689574	15.47%	1.039%
25	10.446	1418	1421	1429	rVV3	72103	114043	2.56%	0.172%
26	10.716	1464	1467	1471	rVV3	50991	57679	1.29%	0.087%
27	11.099	1530	1532	1536	rVB2	43732	41738	0.94%	0.063%
28	11.146	1536	1540	1542	rBV3	29958	48614	1.09%	0.073%
29	11.293	1562	1565	1568	rBV	2652644	2354101	52.80%	3.548%
30	11.316	1568	1569	1572	rVV	937347	652129	14.63%	0.983%
31	11.352	1572	1575	1578	rVV	3996343	3490186	78.28%	5.260%
32	11.404	1581	1584	1588	rBV2	51743	60212	1.35%	0.091%
33	11.528	1600	1605	1607	rBV	100017	100950	2.26%	0.152%
34	11.634	1620	1623	1628	rVV2	61403	74421	1.67%	0.112%

LSC Area Percent Report

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Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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_P\METHODS\SOM-EPA-BP100919MA.M
Title : SVOA CALIBRATION

35	11.716	1634	1637	1641	rVB	37198	41617	0.93%	0.063%
36	11.804	1648	1652	1654	rVV	234513	218726	4.91%	0.330%
37	11.834	1654	1657	1661	rVV2	361236	457324	10.26%	0.689%
38	11.881	1661	1665	1667	rVV2	126749	177624	3.98%	0.268%
39	11.916	1667	1671	1673	rVV2	443454	471860	10.58%	0.711%
40	11.940	1673	1675	1680	rVB	195059	178271	4.00%	0.269%
41	12.040	1690	1692	1695	rVV	46027	43144	0.97%	0.065%
42	12.099	1699	1702	1705	rVV	148675	164459	3.69%	0.248%
43	12.128	1705	1707	1711	rVB2	159552	144772	3.25%	0.218%
44	12.257	1726	1729	1731	rVV	96461	97591	2.19%	0.147%
45	12.287	1731	1734	1736	rVV	99897	88523	1.99%	0.133%
46	12.310	1736	1738	1740	rVB	69351	50476	1.13%	0.076%
47	12.363	1743	1747	1749	rBV	232999	243022	5.45%	0.366%
48	12.399	1749	1753	1755	rVV3	148862	199168	4.47%	0.300%
49	12.446	1758	1761	1766	rVB2	163890	193143	4.33%	0.291%
50	12.522	1770	1774	1778	rVB	2252707	1919378	43.05%	2.893%
51	12.569	1779	1782	1784	rBV2	86484	80131	1.80%	0.121%
52	12.622	1788	1791	1794	rVB2	47135	40792	0.91%	0.061%
53	12.663	1795	1798	1804	rBV5	89399	130865	2.94%	0.197%
54	12.740	1807	1811	1813	rBV	4209594	4458674	100.00%	6.720%
55	12.757	1813	1814	1817	rVV	2268941	1131849	25.39%	1.706%
56	12.781	1817	1818	1820	rVB2	76977	41475	0.93%	0.063%
57	12.804	1820	1822	1827	rVB2	99355	144091	3.23%	0.217%
58	12.875	1831	1834	1836	rBV	120120	102307	2.29%	0.154%
59	12.916	1838	1841	1845	rVB	130842	147943	3.32%	0.223%
60	12.975	1847	1851	1855	rBV2	209457	295675	6.63%	0.446%
61	13.034	1859	1861	1863	rBV2	56831	58958	1.32%	0.089%
62	13.063	1863	1866	1867	rBV3	160430	135737	3.04%	0.205%
63	13.087	1867	1870	1873	rVB	450007	427873	9.60%	0.645%
64	13.122	1873	1876	1878	rBV2	60460	72643	1.63%	0.109%
65	13.151	1878	1881	1884	rVB	307105	313212	7.02%	0.472%
66	13.193	1884	1888	1891	rBV2	345949	346094	7.76%	0.522%
67	13.287	1901	1904	1907	rVV	215213	183328	4.11%	0.276%
68	13.316	1907	1909	1913	rBV	228102	191767	4.30%	0.289%
69	13.381	1918	1920	1924	rBV	160827	150416	3.37%	0.227%
70	13.493	1937	1939	1942	rBV2	106168	101261	2.27%	0.153%
71	13.604	1956	1958	1963	rVB	155455	170271	3.82%	0.257%

LSC Area Percent Report

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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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_P\METHODS\SOM-EPA-BP100919MA.M
Title : SVOA CALIBRATION

72	13.716	1973	1977	1979	rBV2	259080	292486	6.56%	0.441%
73	13.740	1979	1981	1984	rVV	185302	182140	4.09%	0.275%
74	13.769	1984	1986	1990	rVV3	201756	254355	5.70%	0.383%
75	13.816	1991	1994	2003	rVB2	950341	1126042	25.26%	1.697%
76	13.969	2015	2020	2023	rBV2	2479733	3483229	78.12%	5.250%
77	13.998	2023	2025	2028	rVB	1017257	808368	18.13%	1.218%
78	14.057	2033	2035	2036	rBV	73609	58830	1.32%	0.089%
79	14.075	2036	2038	2041	rVB	162092	123323	2.77%	0.186%
80	14.116	2041	2045	2047	rBV3	79374	134815	3.02%	0.203%
81	14.145	2047	2050	2057	rBV6	256382	257618	5.78%	0.388%
82	14.328	2079	2081	2083	rBV	96829	87277	1.96%	0.132%
83	14.363	2083	2087	2090	rVV	345047	391440	8.78%	0.590%
84	14.398	2090	2093	2096	rVV	200363	207801	4.66%	0.313%
85	14.457	2100	2103	2106	rVV2	376758	433911	9.73%	0.654%
86	14.493	2106	2109	2111	rVV	207174	254524	5.71%	0.384%
87	14.510	2111	2112	2116	rVB	197716	137922	3.09%	0.208%
88	14.769	2152	2156	2158	rBV2	253356	283789	6.36%	0.428%
89	14.840	2165	2168	2171	rBV3	122585	163767	3.67%	0.247%
90	15.022	2194	2199	2207	rBV	999583	1858483	41.68%	2.801%
91	15.104	2210	2213	2216	rBV2	246504	323651	7.26%	0.488%
92	15.322	2246	2250	2253	rBV	517324	727845	16.32%	1.097%
93	15.363	2253	2257	2259	rVV	2350798	2919081	65.47%	4.399%
94	15.387	2259	2261	2267	rVB	902362	924891	20.74%	1.394%
95	15.451	2267	2272	2275	rBV	1479649	1805864	40.50%	2.722%
96	15.487	2275	2278	2284	rVB3	326113	493643	11.07%	0.744%
97	15.928	2349	2353	2358	rBV2	191898	278499	6.25%	0.420%
98	16.439	2436	2440	2444	rBV	90744	146444	3.28%	0.221%
99	16.916	2516	2521	2531	rVB2	363528	715638	16.05%	1.079%
100	17.357	2591	2596	2609	rVB	239998	472924	10.61%	0.713%

Sum of corrected areas: 66350707

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000762.D
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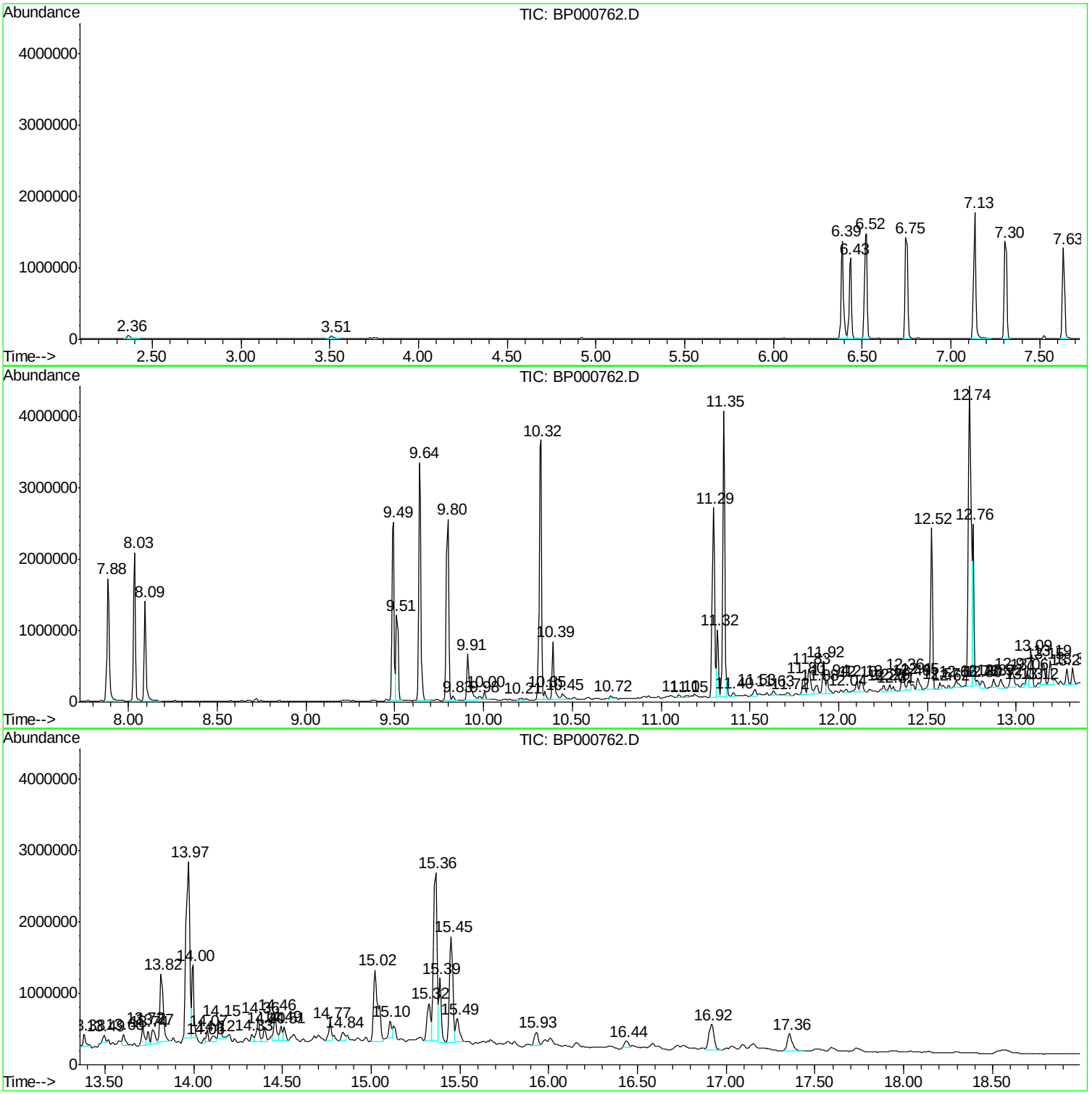
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BEPF8

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

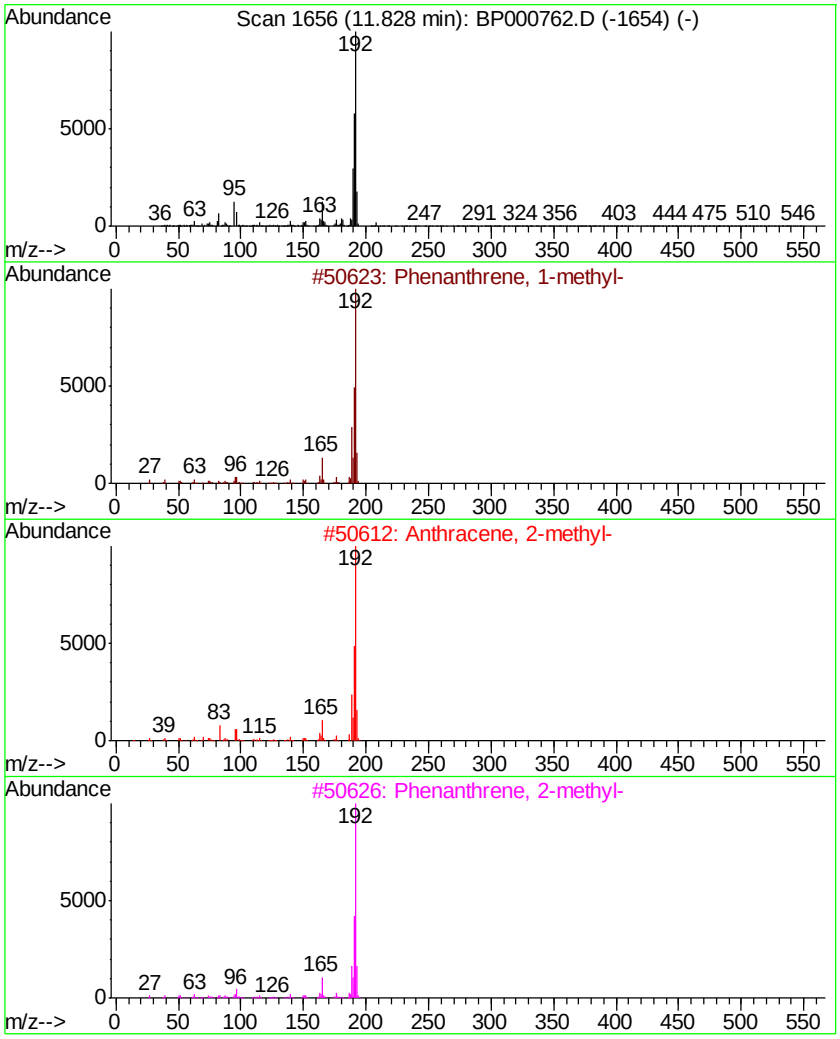
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

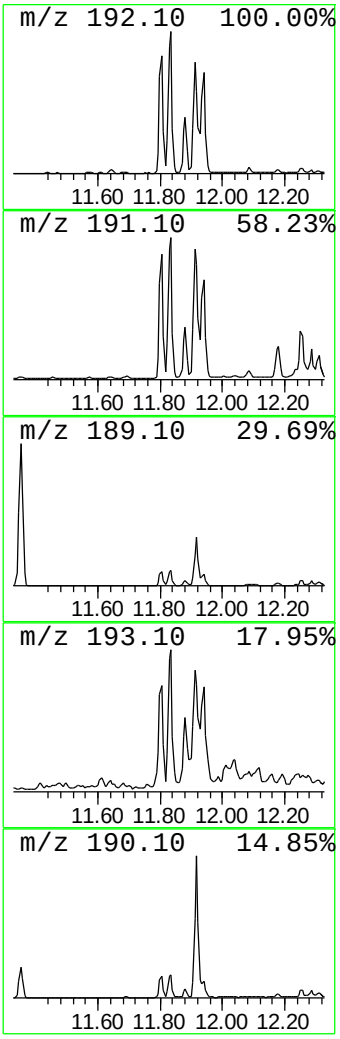
****

Peak Number 1 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	3.89 ng/ul	457324	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101619\
Data File : BP000762.D
Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

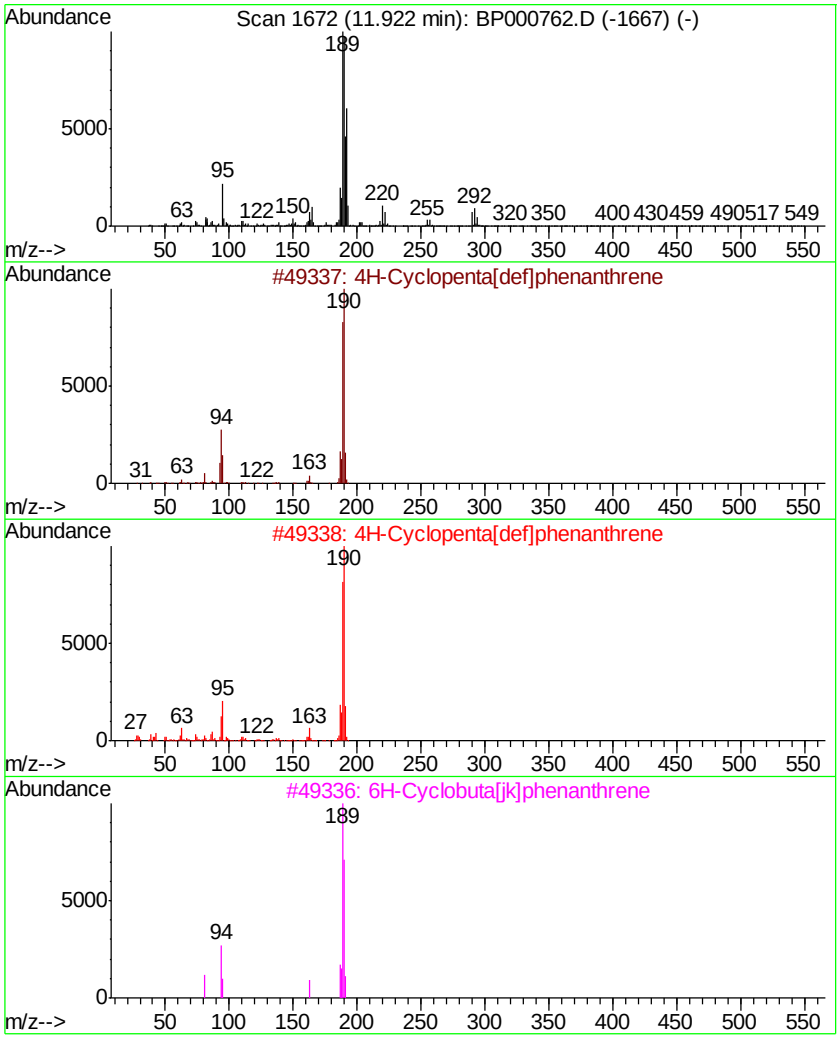
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

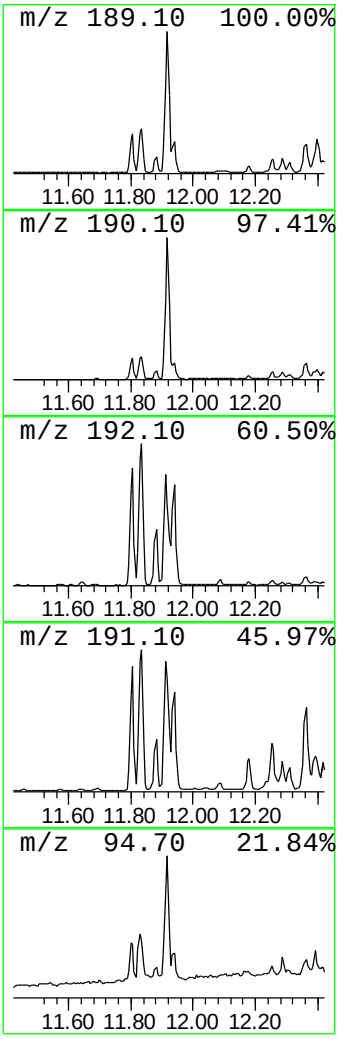
****

Peak Number 2 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	4.01 ng/ul	471860	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
4	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	47





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Data File : BP000762.D
Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

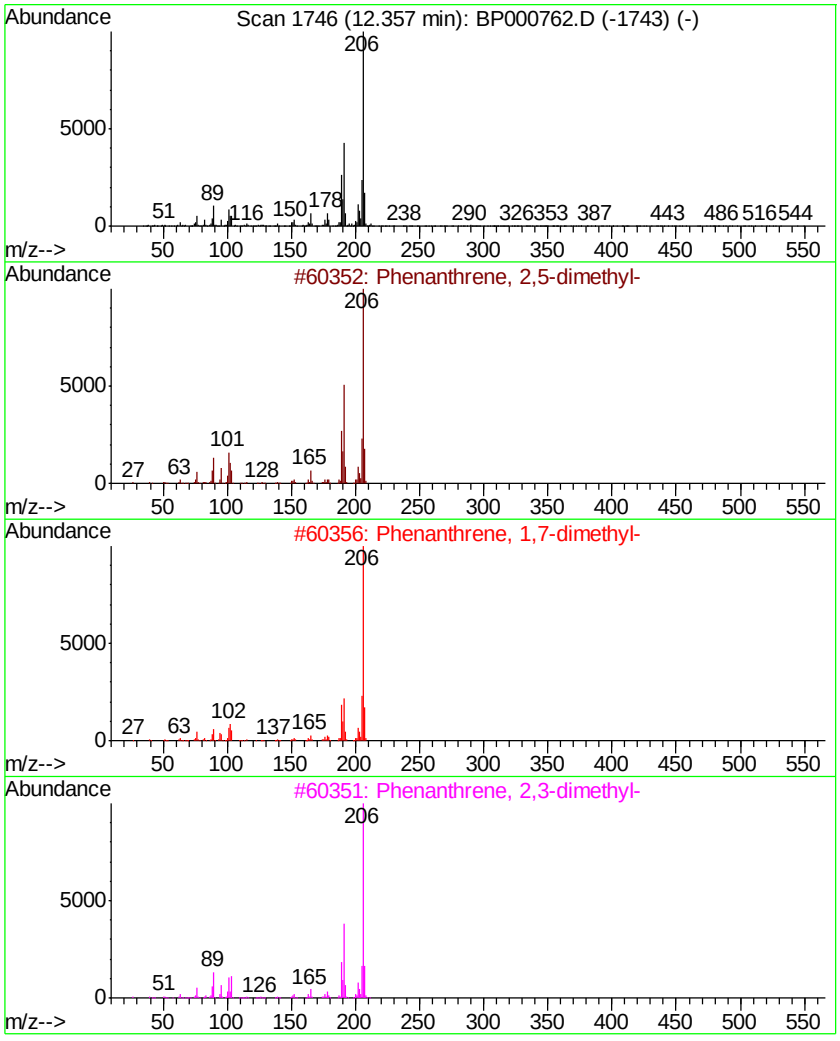
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

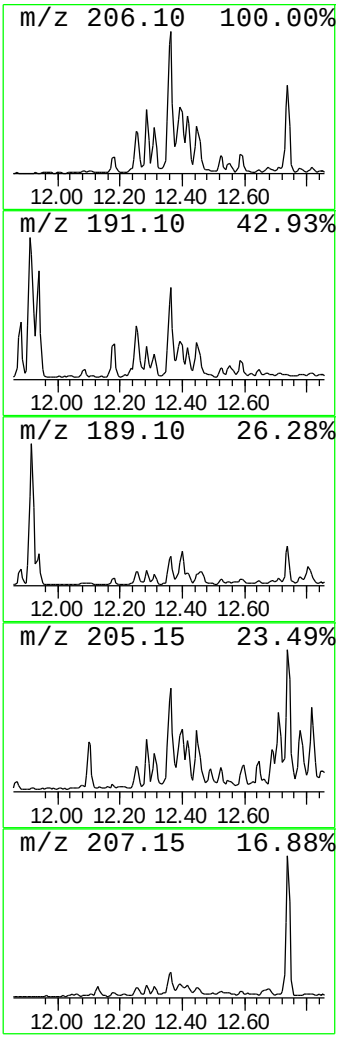
****

Peak Number 3 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	2.06 ng/ul	243022	Phenanthrene-d10	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4	di-p-Tolylacetylene	206	C16H14	002789-88-0	86
5	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	81





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Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

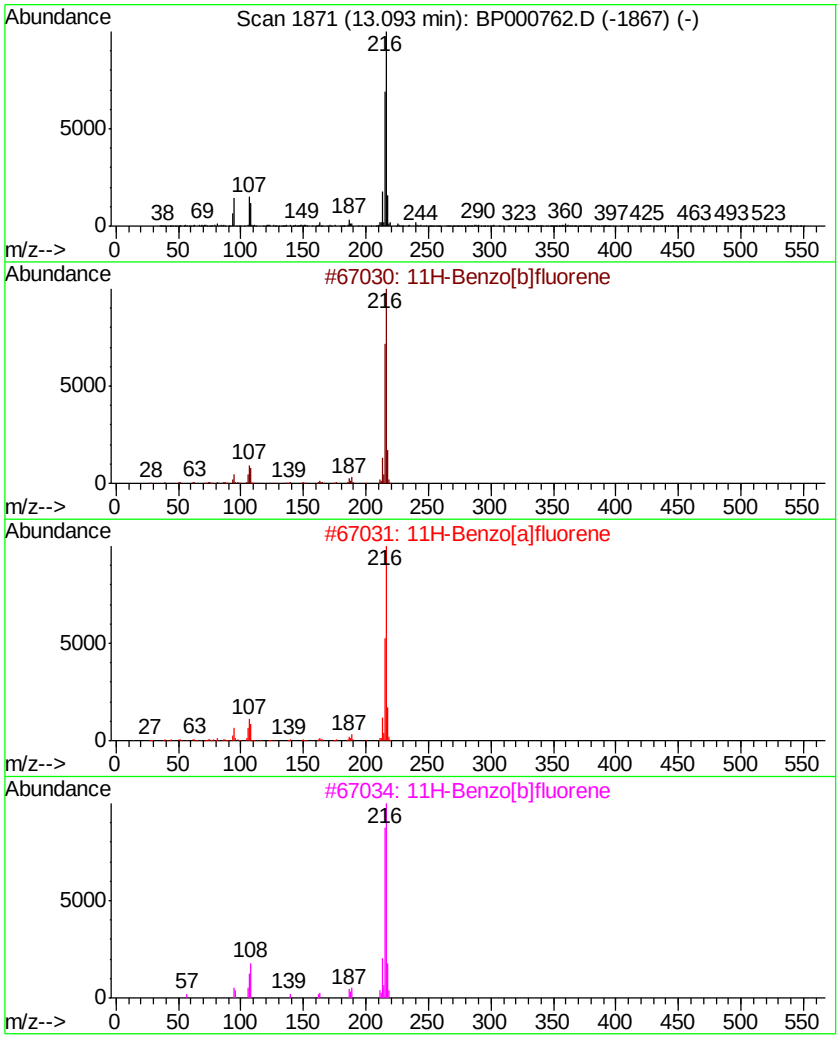
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

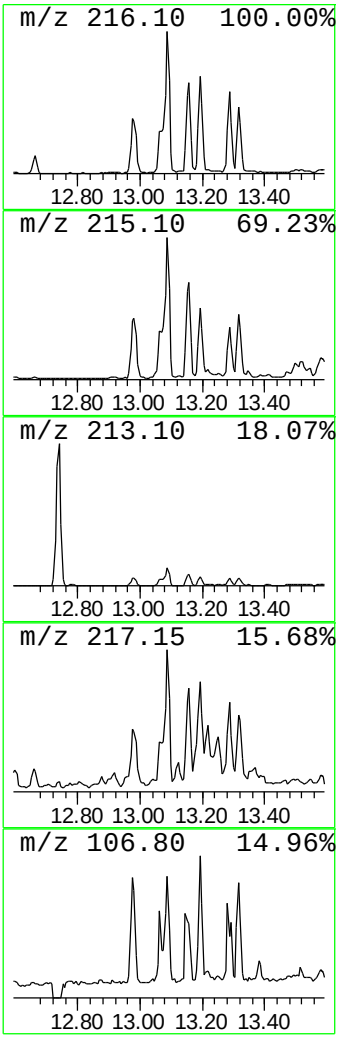
****

Peak Number 4 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	2.46 ng/ul	427873	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	72





Library Search Compound Report

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Data File : BP000762.D
Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

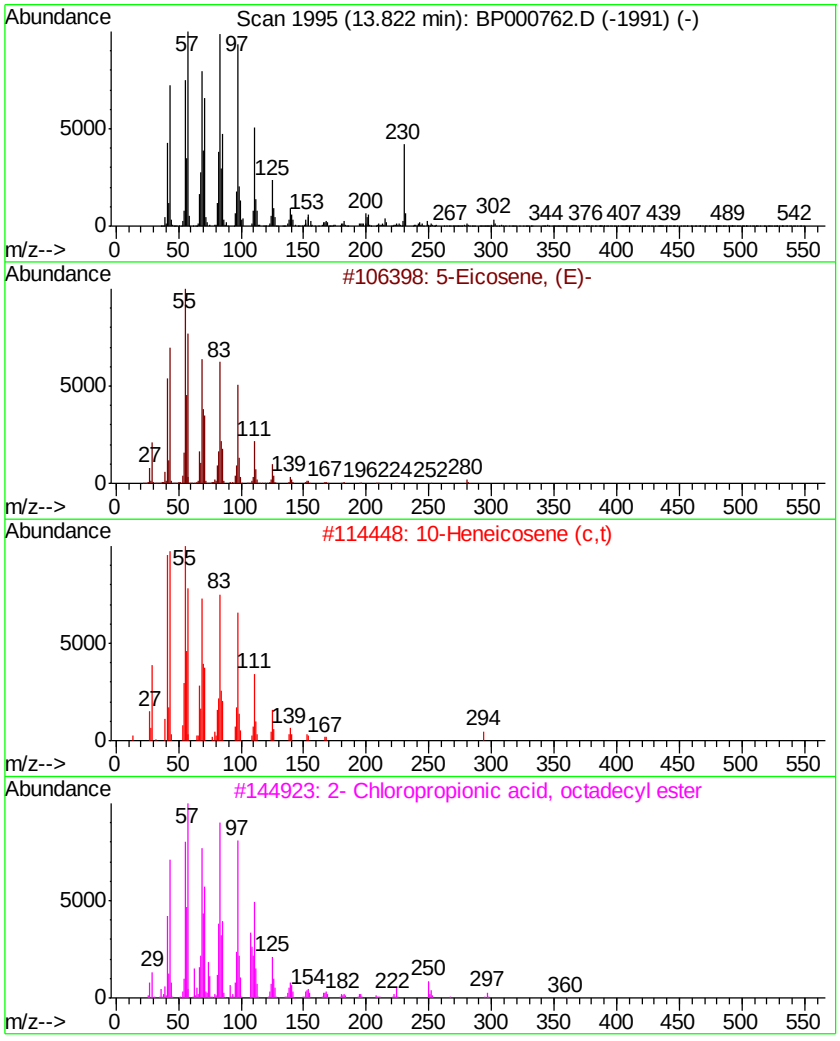
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

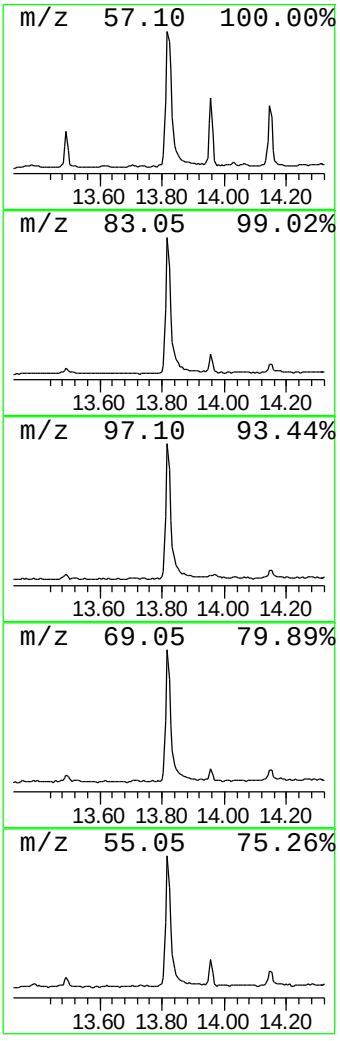
****

Peak Number 5 (DEL) Alkane: Cyclic13.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	6.47 ng/ul	1126040	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	10-Heneicosene (c,t)	294	C21H42	095008-11-0	94
3	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
4	1-Nonadecene	266	C19H38	018435-45-5	93
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	90





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Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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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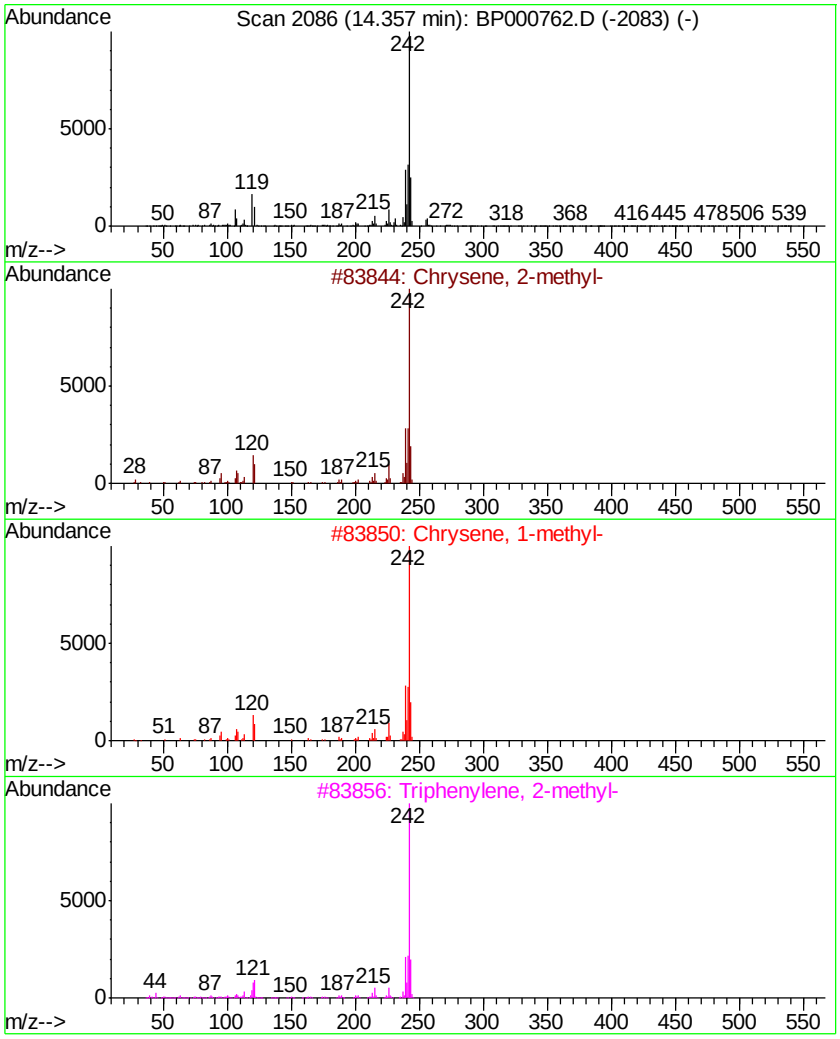
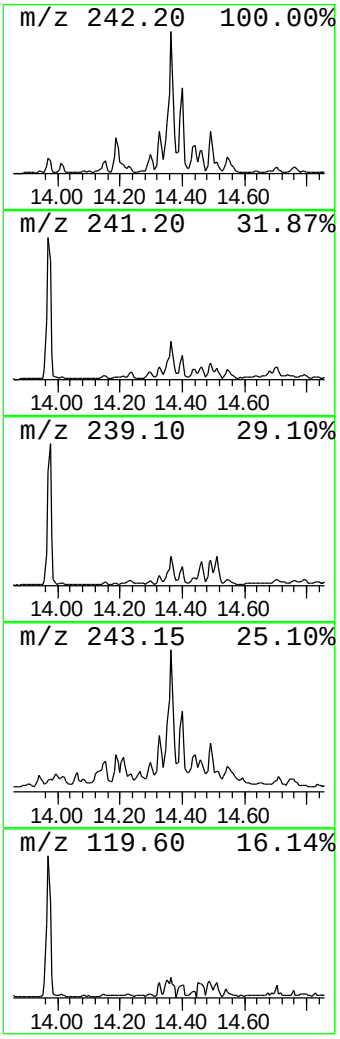
Peak Number 6 Chrysene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.36	2.25 ng/ul	391440	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chrysene, 2-methyl-	242	C19H14	003351-32-4	98
2	Chrysene, 1-methyl-	242	C19H14	003351-28-8	97
3	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	97
4	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	95
5	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	95





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Acq On : 15 Oct 2019 13:45
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Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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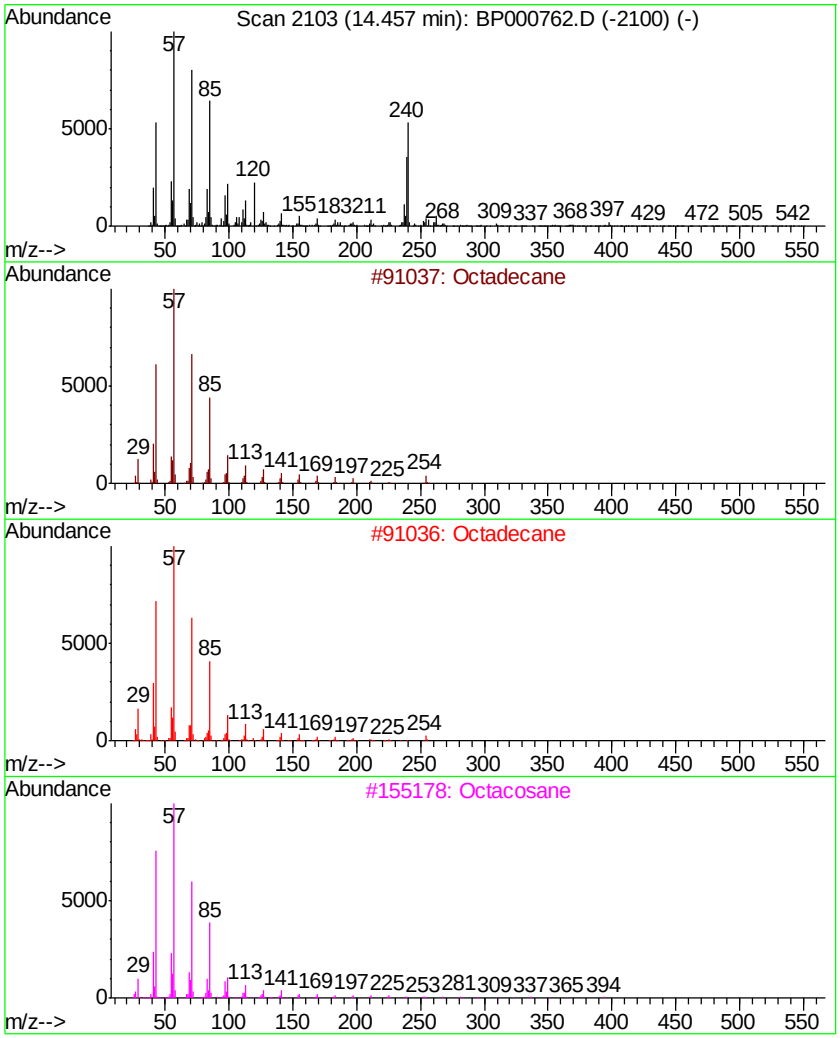
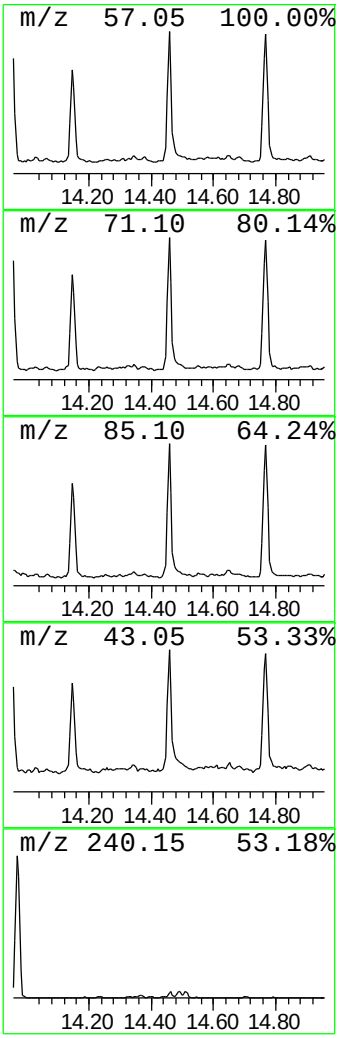
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	2.49 ng/ul	433911	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	95
2	Octadecane	254	C18H38	000593-45-3	91
3	Octacosane	394	C28H58	000630-02-4	86
4	Octadecane	254	C18H38	000593-45-3	70
5	Heneicosane	296	C21H44	000629-94-7	55





Library Search Compound Report

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Acq On : 15 Oct 2019 13:45
Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

TIC Library : C:\DATABASE\NIST02.L
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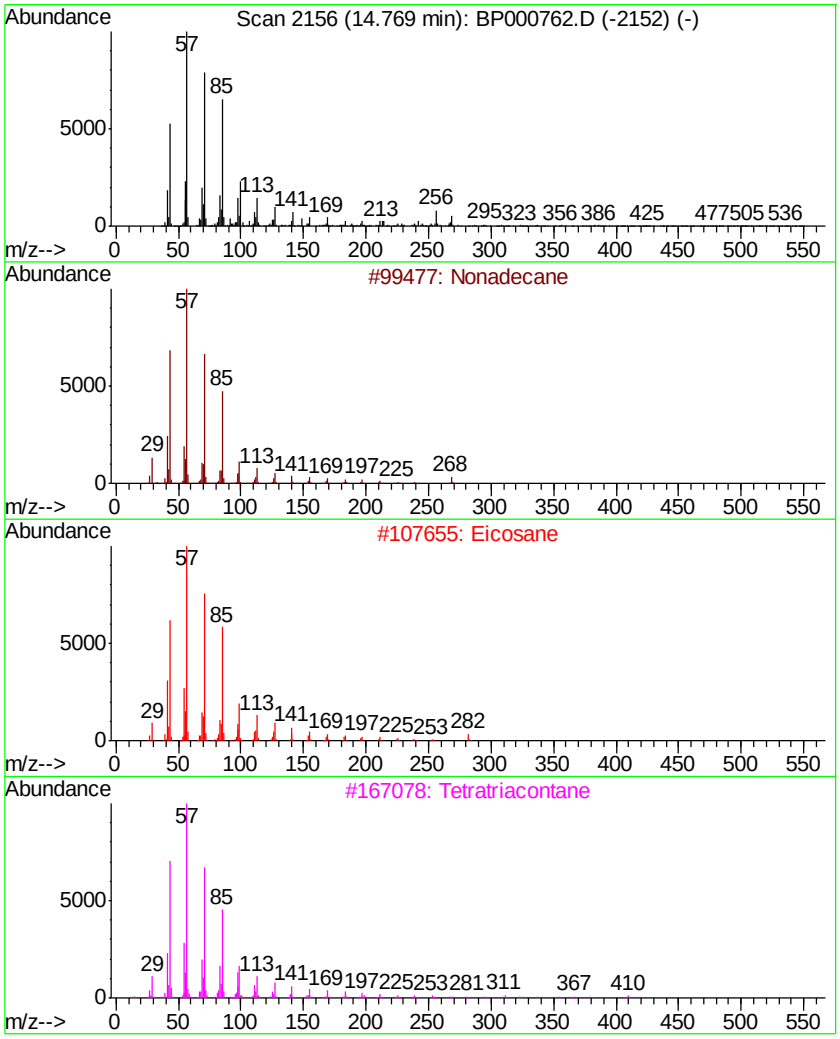
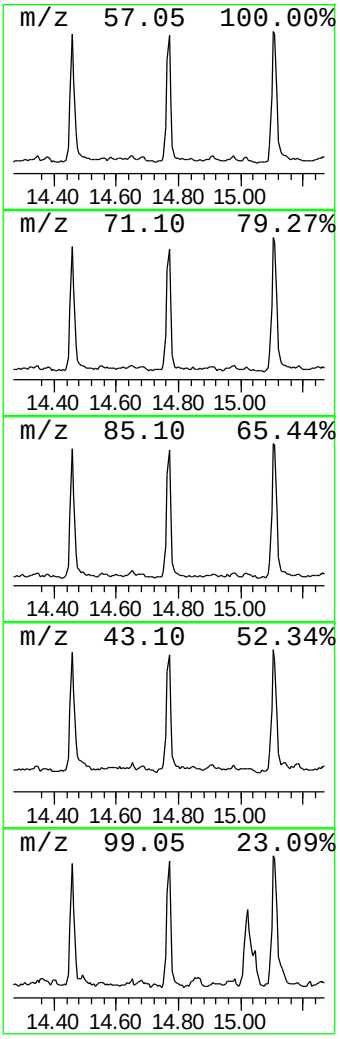
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	3.14 ng/ul	283789	Perylene-d12	15.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	96
2	Eicosane	282	C20H42	000112-95-8	90
3	Tetratriacontane	479	C34H70	014167-59-0	87
4	Heptadecane	240	C17H36	000629-78-7	87
5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87





Library Search Compound Report

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Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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Quant Title : SVOA CALIBRATION

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

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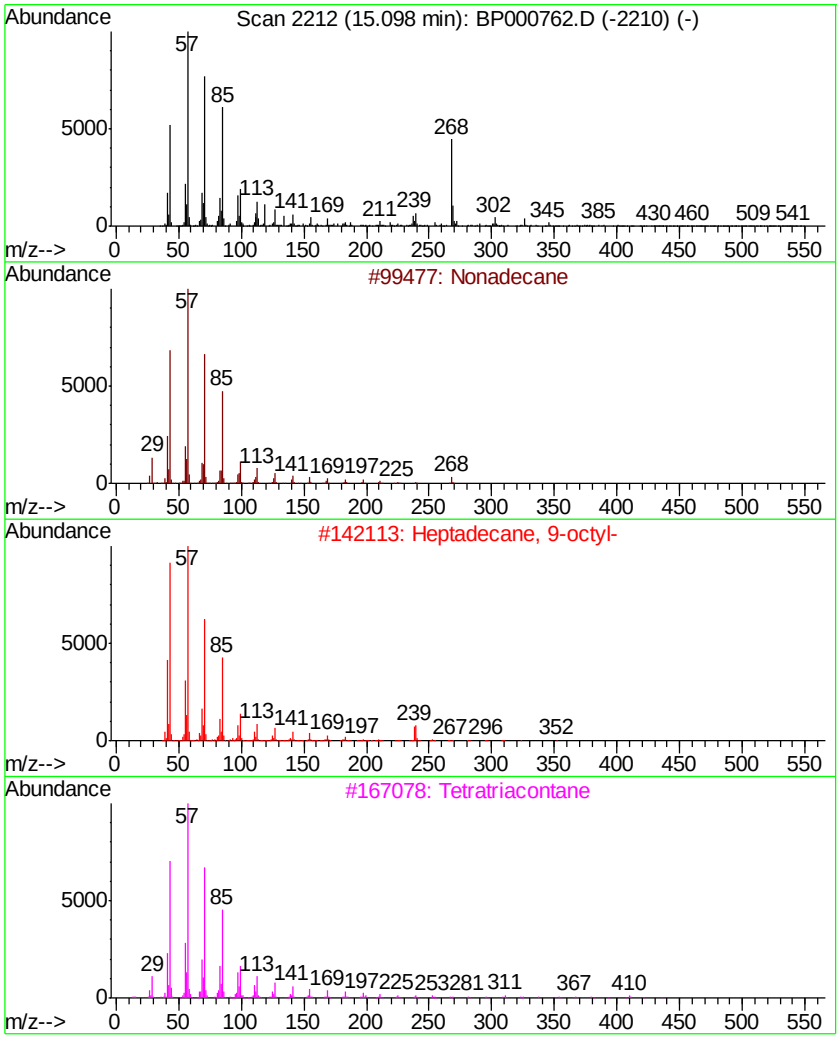
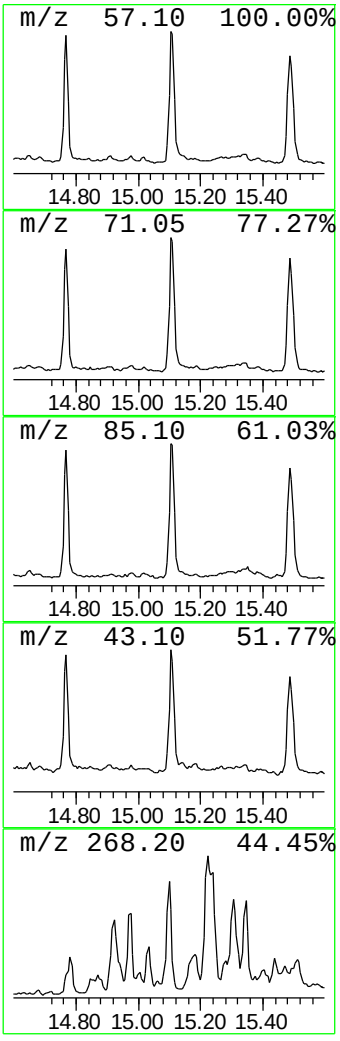
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	3.58 ng/ul	323651	Perylene-d12	15.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	93
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3	Tetratriacontane	479	C34H70	014167-59-0	90
4	Heneicosane	296	C21H44	000629-94-7	90
5	Triacontane	422	C30H62	000638-68-6	87





Library Search Compound Report

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Operator : JU
Sample : K5175-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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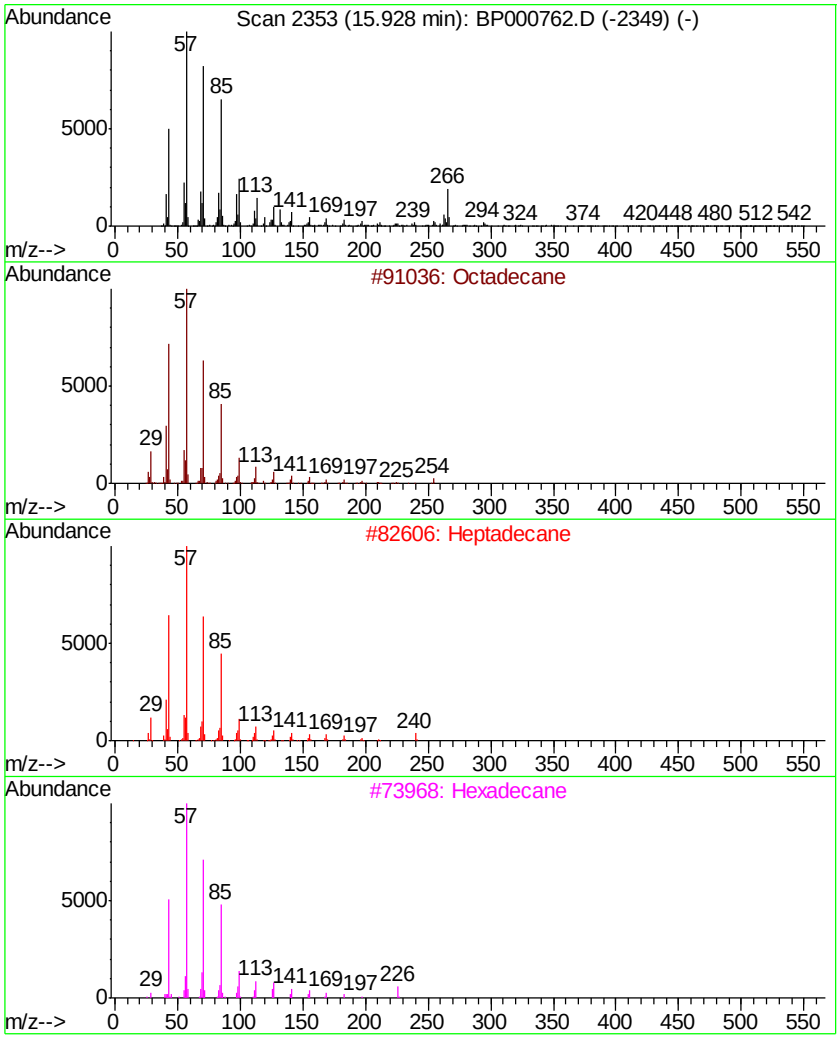
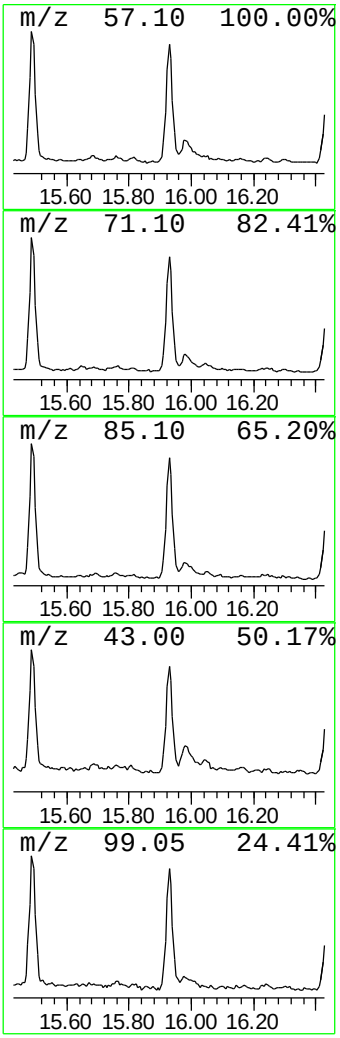
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.93	3.08 ng/ul	278499	Perylene-d12	15.45

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	96
2	Heptadecane	240	C17H36	000629-78-7	90
3	Hexadecane	226	C16H34	000544-76-3	89
4	Heneicosane	296	C21H44	000629-94-7	87
5	Triacontane	422	C30H62	000638-68-6	87





Tentatively Identified Compound (LSC) summary

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Data File : BP000762.D
Acq On : 15 Oct 2019 13:45
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : BEPF8

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

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Phenanthrene, 1-m...	11.83	3.9	ng/ul	457324	4	11.29	2354100	20.0
4H-Cyclopenta[def...	11.92	4.0	ng/ul	471860	4	11.29	2354100	20.0
Phenanthrene, 2,5...	12.36	2.1	ng/ul	243022	4	11.29	2354100	20.0
11H-Benzo[b]fluorene	13.09	2.5	ng/ul	427873	5	13.97	3483230	20.0
(DEL) Alkane: Cyc...	13.82	6.5	ng/ul	1126040	5	13.97	3483230	20.0
Chrysene, 2-methyl-	14.36	2.3	ng/ul	391440	5	13.97	3483230	20.0
(DEL) Alkane: Str...	14.46	2.5	ng/ul	433911	5	13.97	3483230	20.0
(DEL) Alkane: Str...	14.77	3.1	ng/ul	283789	6	15.45	1805860	20.0
(DEL) Alkane: Str...	15.10	3.6	ng/ul	323651	6	15.45	1805860	20.0
(DEL) Alkane: Str...	15.93	3.1	ng/ul	278499	6	15.45	1805860	20.0