LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBXZ2

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\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

Signal : TIC: BP013158.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	22	25	32	rVB	132743	167630	1.26%	0.108%
2	3.764	95	98	113	rBV	977384	1484397	11.16%	0.956%
3	3.870	113	116	125	rVB	62064	87544	0.66%	0.056%
4	3.999	134	138	142	rVB	45204	56532	0.42%	0.036%
5	4.093	151	154	158	rBV	33012	40069	0.30%	0.026%
6	4.323	189	193	202	rVB	86743	114364	0.86%	0.074%
7	4.664	248	251	255	rVB5	14709	15721	0.12%	0.010%
8	4.740	260	264	269	rVB	53962	72002	0.54%	0.046%
9	5.040	309	315	323	rBV	255618	368614	2.77%	0.237%
10	6.628	579	585	589	rVB9	10498	20515	0.15%	0.013%
11	6.928	631	636	641	rVV8	9585	19755	0.15%	0.013%
12	7.111	657	667	680	rBV2	2158310	4354211	32.73%	2.805%
13	7.281	689	696	707	rVB	1684925	2594180	19.50%	1.671%
14	7.481	723	730	740	rBV	2098867	3218779	24.19%	2.073%
15	7.952	802	810	817	rBV	1565611	2433680	18.29%	1.568%
16	8.016	817	821	827	rVB6	11927	20994	0.16%	0.014%
17	8.663	923	931	940	rBV	2261600	3592868	27.01%	2.314%
18	8.781	946	951	958	rVB	51599	92273	0.69%	0.059%
19	9.122	1002	1009	1018	rBV	1941706	3110238	23.38%	2.004%
20	9.522	1073	1077	1083	rVB2	35901	54151	0.41%	0.035%
21	9.846	1122	1132	1144	rBV	1598725	2626852	19.75%	1.692%
22	10.069	1166	1170	1176	rVB8	11534	20934	0.16%	0.013%
23	10.387	1212	1224	1239	rBV	2373694	4008644	30.13%	2.582%
24	10.769	1278	1289	1303	rBV	2039238	3444550	25.89%	2.219%
25	10.905	1305	1312	1326	rBV	1525617	2744349	20.63%	1.768%
26	11.552	1416	1422	1430	rBV9	12033	28628	0.22%	0.018%
27	11.681	1439	1444	1449	rBV9	6543	13434	0.10%	0.009%
28	11.787	1451	1462	1467	rBV9	10555	26218	0.20%	0.017%
29	12.087	1507	1513	1517	rBV4	12435	17816	0.13%	0.011%
30	12.175	1523	1528	1533	rBV	34385	45024	0.34%	0.029%
31	12.293	1535	1548	1553	rBV3	24553	56911	0.43%	0.037%
32	12.352	1553	1558	1564	rVV2	43727	75381	0.57%	0.049%
33	12.457	1567	1576	1588	rVV	436308	788683	5.93%	0.508%
34	12.557	1591	1593	1599	rVV7	11759	22304	0.17%	0.014%
35	12.646	1602	1608	1614	rVB8	10226	21528	0.16%	0.014%
36	12.863	1639	1645	1646	rBV6	10826	14358	0.11%	0.009%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
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Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

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\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

37	13.128	1685	1690	1693	rBV6	12173	18218	0.14%	0.012%
38	13.410	1727	1738	1744	rVB3	59851	100850	0.76%	0.065%
39	13.487	1744	1751	1753	rBV8	11734	26684	0.20%	0.017%
40	13.563	1761	1764	1770	rVB8	8879	18514	0.14%	0.012%
41	13.916	1819	1824	1830	rBV6	57352	85096	0.64%	0.055%
42	13.998	1830	1838	1848	rBV	3876319	5550621	41.72%	3.576%
43	14.063	1848	1849	1854	rVB4	18678	17275	0.13%	0.011%
44	14.116	1854	1858	1865	rVB9	20176	34087	0.26%	0.022%
45	14.181	1865	1869	1874	rBV	48438	64270	0.48%	0.041%
46	14.287	1881	1887	1897	rBV	4591607	6829904	51.34%	4.400%
47	14.481	1915	1920	1925	rVB2	38828	57801	0.43%	0.037%
48	14.593	1933	1939	1947	rVB	2821795	4318191	32.46%	2.782%
49	14.787	1965	1972	1987	rBV	1415743	2314631	17.40%	1.491%
50	14.946	1995	1999	2004	rVV6	21685	31990	0.24%	0.021%
51	15.010	2004	2010	2023	rVB7	49544	111025	0.83%	0.072%
52	15.381	2068	2073	2075	rBV6	13972	18580	0.14%	0.012%
53	15.434	2079	2082	2087	rVB4	16052	21250	0.16%	0.014%
54	15.587	2101	2108	2115	rBV	5760790	8402039	63.16%	5.412%
55	15.704	2122	2128	2140	rBV	1699705	2321311	17.45%	1.495%
56	15.828	2145	2149	2154	rVB4	27410	44673	0.34%	0.029%
57	15.957	2165	2171	2183	rVB	2741620	3723002	27.99%	2.398%
58	16.163	2202	2206	2213	rVB9	13168	24148	0.18%	0.016%
59	16.304	2225	2230	2231	rBV5	20591	31817	0.24%	0.020%
60	16.440	2247	2253	2257	rBV6	28453	46144	0.35%	0.030%
61	16.522	2262	2267	2274	rVB	213238	290315	2.18%	0.187%
62	16.663	2286	2291	2296	rVB6	8685	17215	0.13%	0.011%
63	16.734	2299	2303	2307	rBV7	27847	36102	0.27%	0.023%
64	17.098	2362	2365	2367	rVB3	12552	14461	0.11%	0.009%
65	17.234	2384	2388	2392	rBV2	61240	85803	0.64%	0.055%
66	17.298	2395	2399	2402	rBV5	33355	46500	0.35%	0.030%
67	17.351	2402	2408	2414	rBV	3429605	4907637	36.89%	3.161%
68	17.451	2419	2425	2434	rVV2	6102734	8526791	64.09%	5.493%
69	17.539	2437	2440	2449	rVB7	48320	87011	0.65%	0.056%
70	18.110	2531	2537	2541	rBV3	60451	111211	0.84%	0.072%
71	18.163	2542	2546	2551	rBV	173638	227885	1.71%	0.147%
72	18.222	2551	2556	2566	rVV	729095	1071335	8.05%	0.690%
73	18.469	2595	2598	2601	rBV	31219	36009	0.27%	0.023%
74	18.510	2601	2605	2610	rBV5	51586	76448	0.57%	0.049%
75	18.786	2649	2652	2657	rVB4	66052	82058	0.62%	0.053%
76	19.122	2703	2709	2714	rBV5	80112	187753	1.41%	0.121%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

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\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

77	19.257	2729	2732	2738	rBV2	92539	115407	0.87%	0.074%
78	19.334	2738	2745	2747	rBV3	216231	419898	3.16%	0.270%
79	19.357	2747	2749	2761	rVB	225989	310601	2.33%	0.200%
80	19.451	2761	2765	2769	rBV2	245145	304542	2.29%	0.196%
81	19.681	2799	2804	2819	rVB	8118508	10856910	81.61%	6.994%
82	19.833	2825	2830	2836	rBV	1022901	1285175	9.66%	0.828%
83	20.192	2885	2891	2895	rBV4	202722	545136	4.10%	0.351%
84	21.128	3046	3050	3058	rBV	1025298	1500204	11.28%	0.966%
85	21.439	3097	3103	3108	rBV	4291335	5859236	44.04%	3.774%
86	23.133	3386	3391	3397	rBV	2562442	4075694	30.64%	2.625%
87	23.757	3490	3497	3513	rVB2	6162782	13303530	100.00%	8.570%
88	23.916	3518	3524	3533	rVB2	3065269	6259661	47.05%	4.032%
89	24.433	3604	3612	3622	rVB2	5738316	12144258	91.29%	7.823%
90	24.539	3623	3630	3641	rVB	2572266	5788089	43.51%	3.728%
91	24.721	3657	3661	3672	rVB8	699646	1541431	11.59%	0.993%
92	26.451	3948	3955	3975	rVB	1479613	5039133	37.88%	3.246%

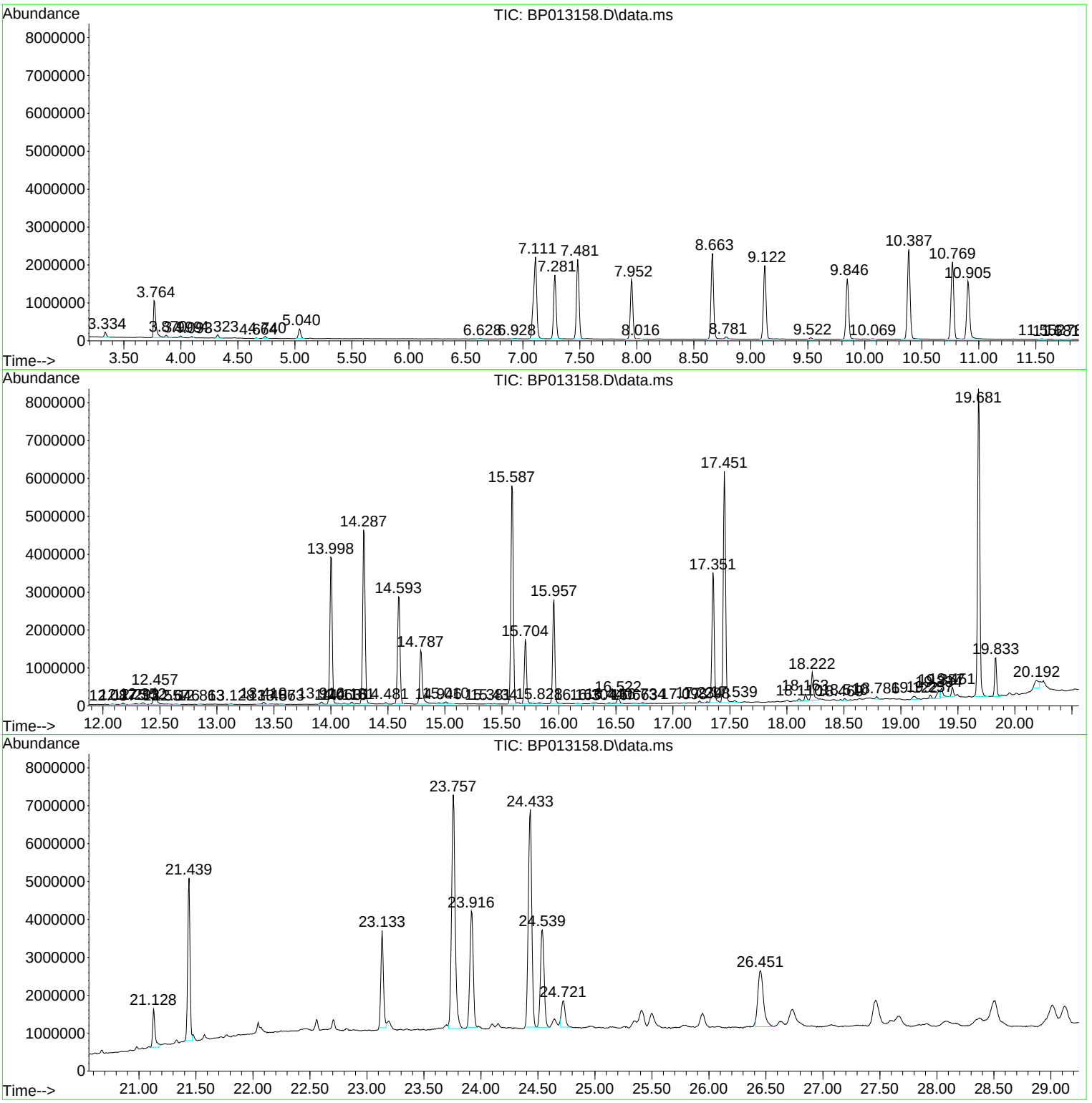
Sum of corrected areas: 155239691

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

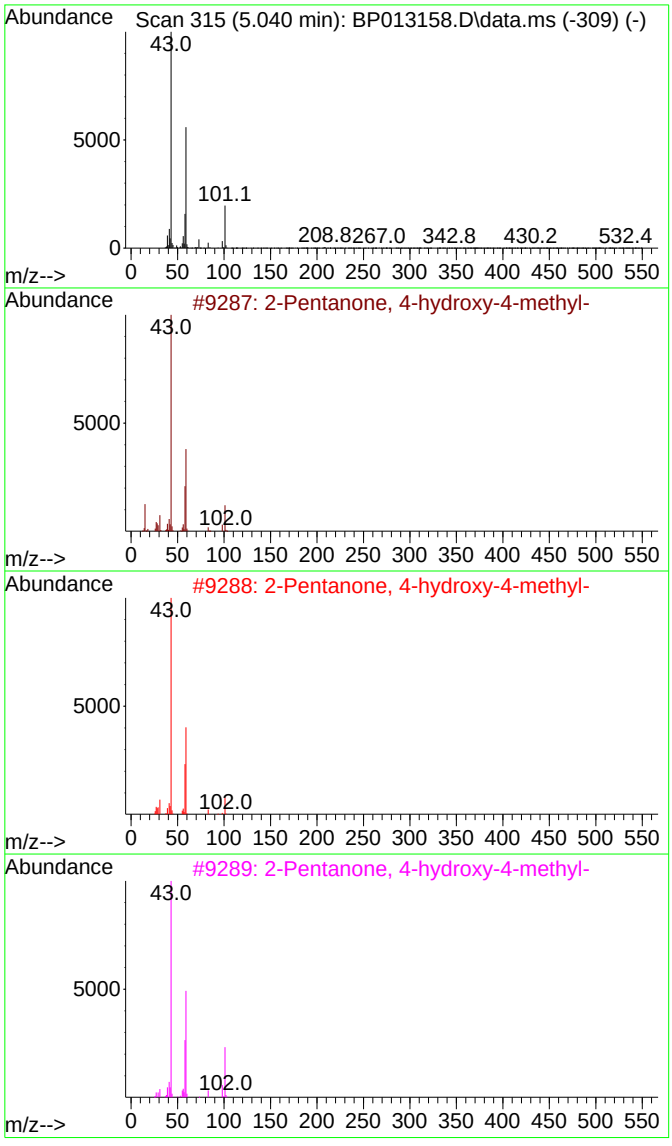
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

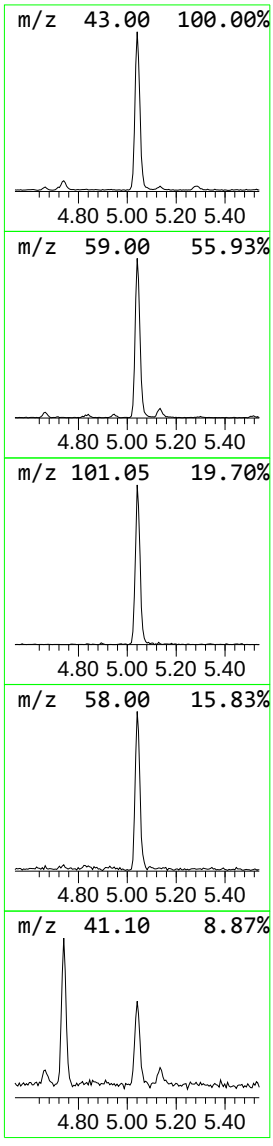
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	3.03 ng/ul	368614	1,4-Dichlorobenzene-d4	7.952

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
4	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	33
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

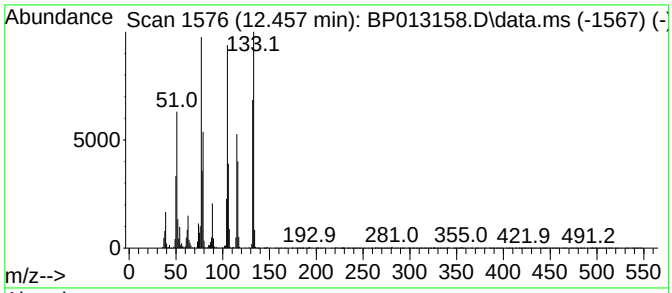
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

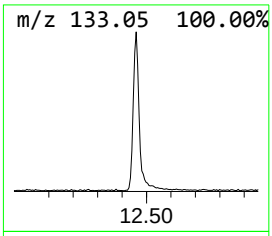
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

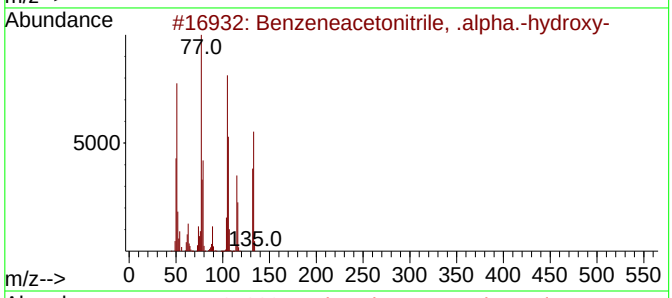
****
Peak Number 2 Benzeneacetonitrile, .alpha... Concentration Rank 8

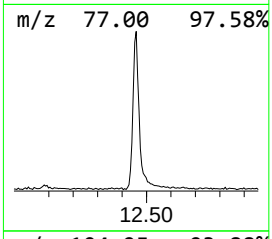
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.457	4.58 ng/ul	788683	Naphthalene-d8	10.769

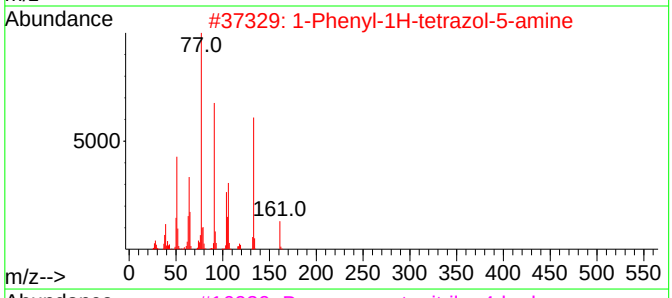
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzeneacetonitrile, .alpha.-hyd...	133	C8H7NO	000532-28-5	87
2	1-Phenyl-1H-tetrazol-5-amine	161	C7H7N5	005467-78-7	47
3	Benzeneacetonitrile, 4-hydroxy-	133	C8H7NO	014191-95-8	38
4	Benzyl isocyanate	133	C8H7NO	003173-56-6	38
5	3-(4-Pyridyl)acrylaldehyde	133	C8H7NO	026505-36-2	35

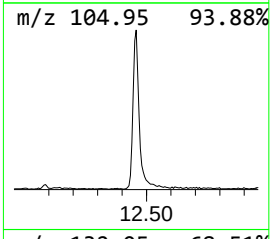


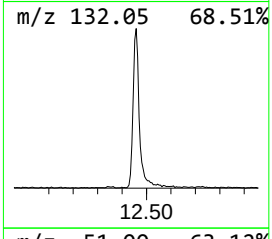


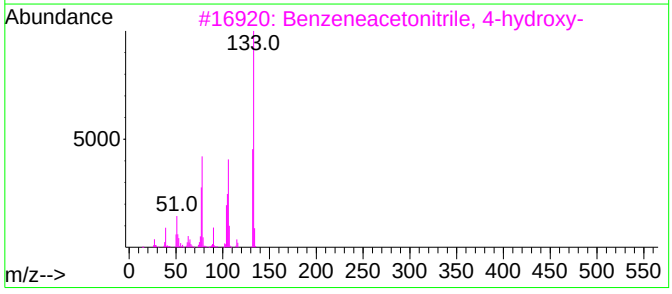


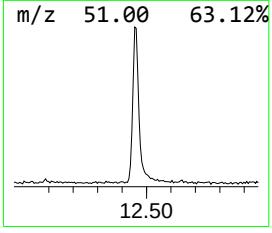












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
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Misc :
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

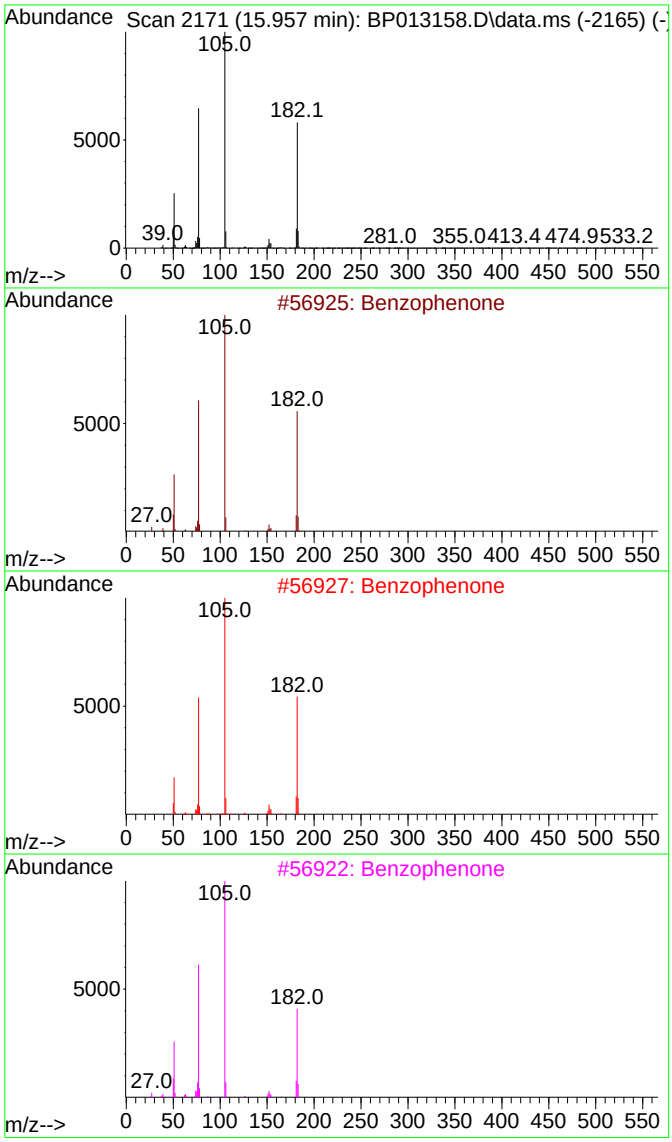
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

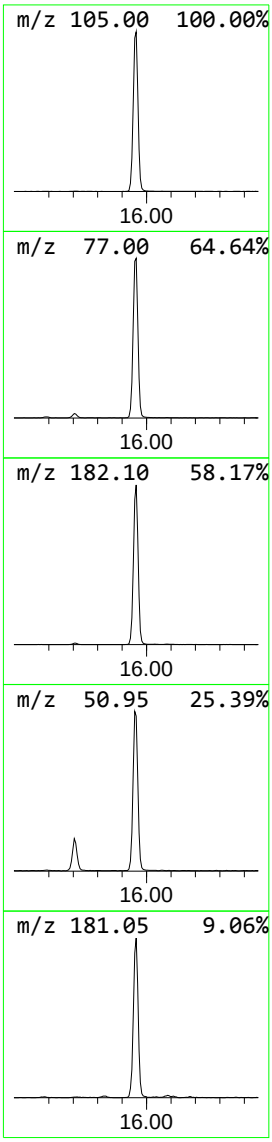
****

Peak Number 3 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.957	17.24 ng/ul	3723000	Acenaphthene-d10	14.593

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	93
5	Benzophenone	182	C13H10O	000119-61-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121922\
Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

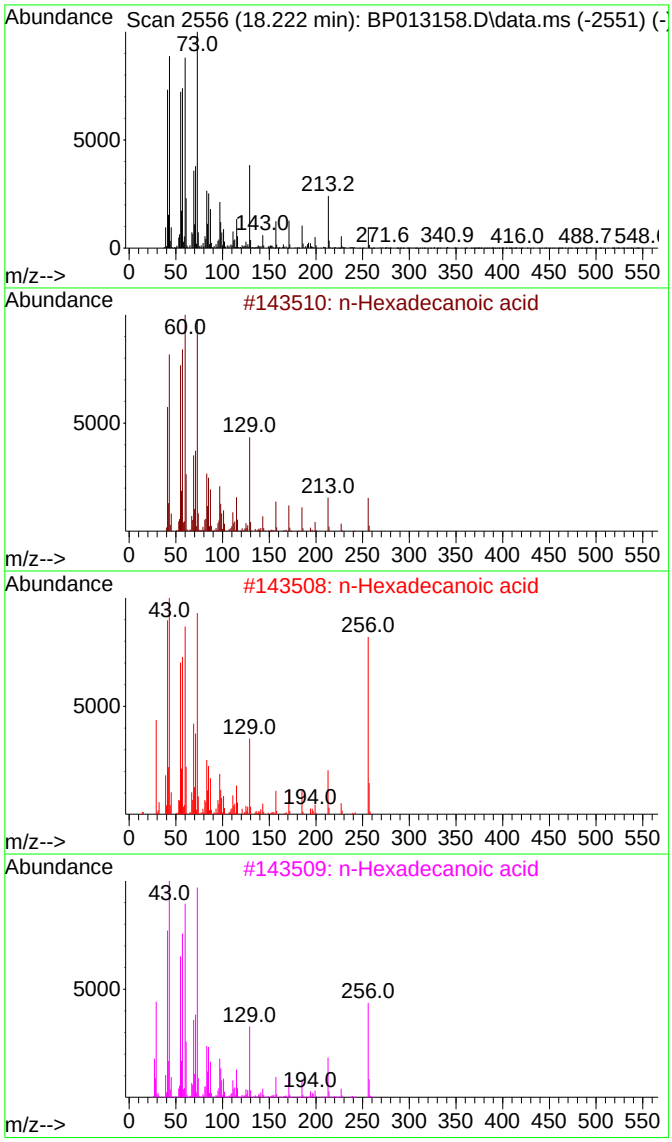
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

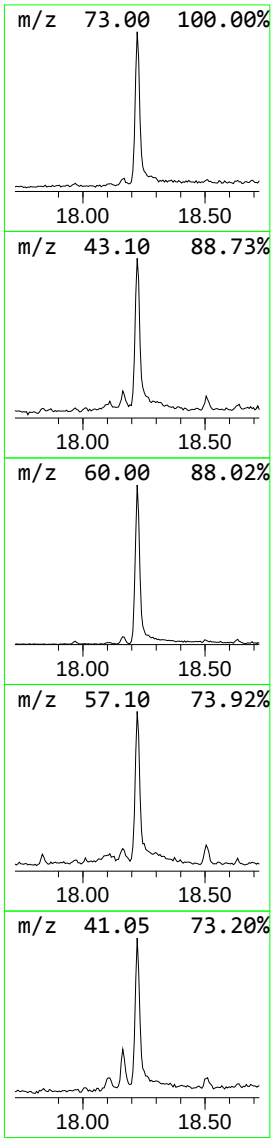
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	4.37 ng/ul	1071340	Phenanthrene-d10	17.351

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	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	93





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Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

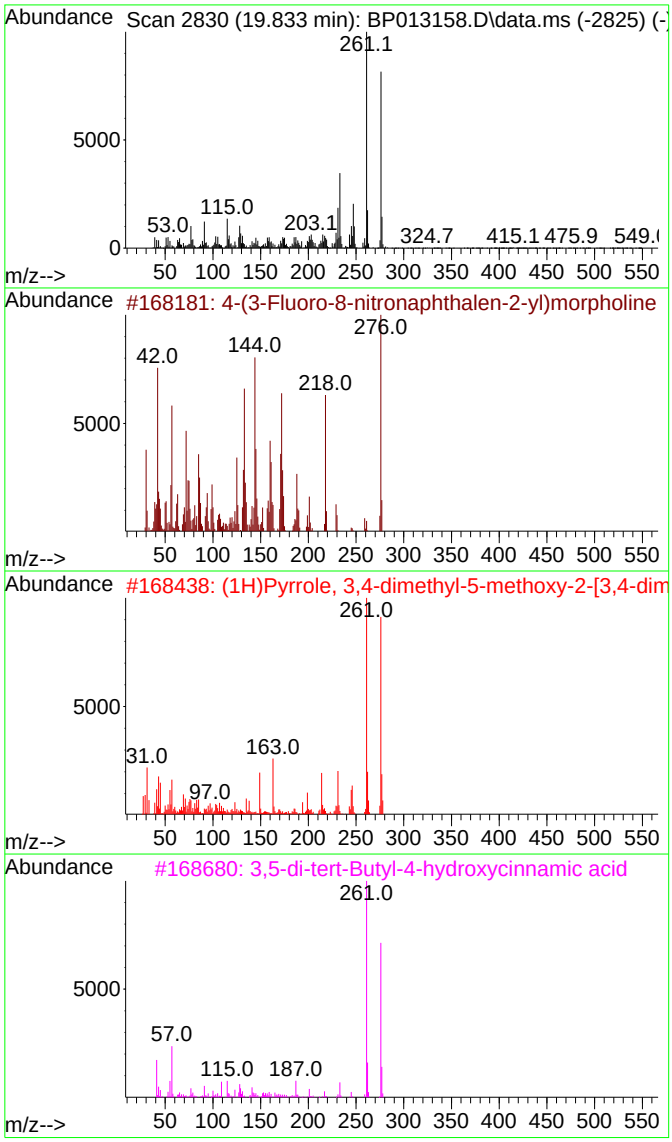
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

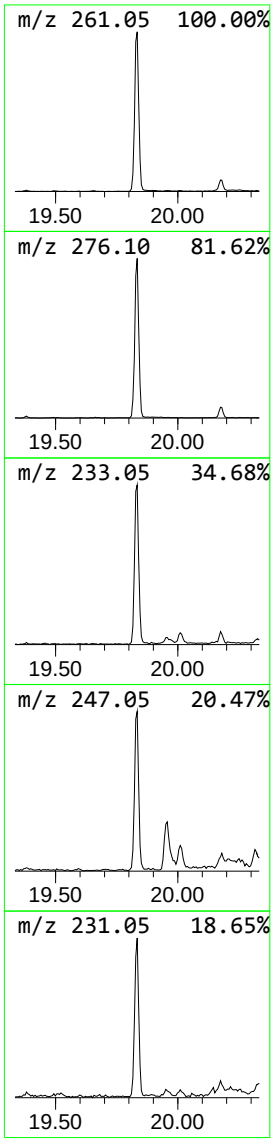
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 4-(3-Fluoro-8-nitronaphthal... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.834	4.39 ng/ul	1285180	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(3-Fluoro-8-nitronaphthalen-2-...	276	C14H13FN2O3	1000409-37-1	92
2	(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	89
3	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	76
4	5-[4-(Diethylamino)phenyl]-4-eth...	276	C14H20N4S	1000476-62-2	72
5	5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	72





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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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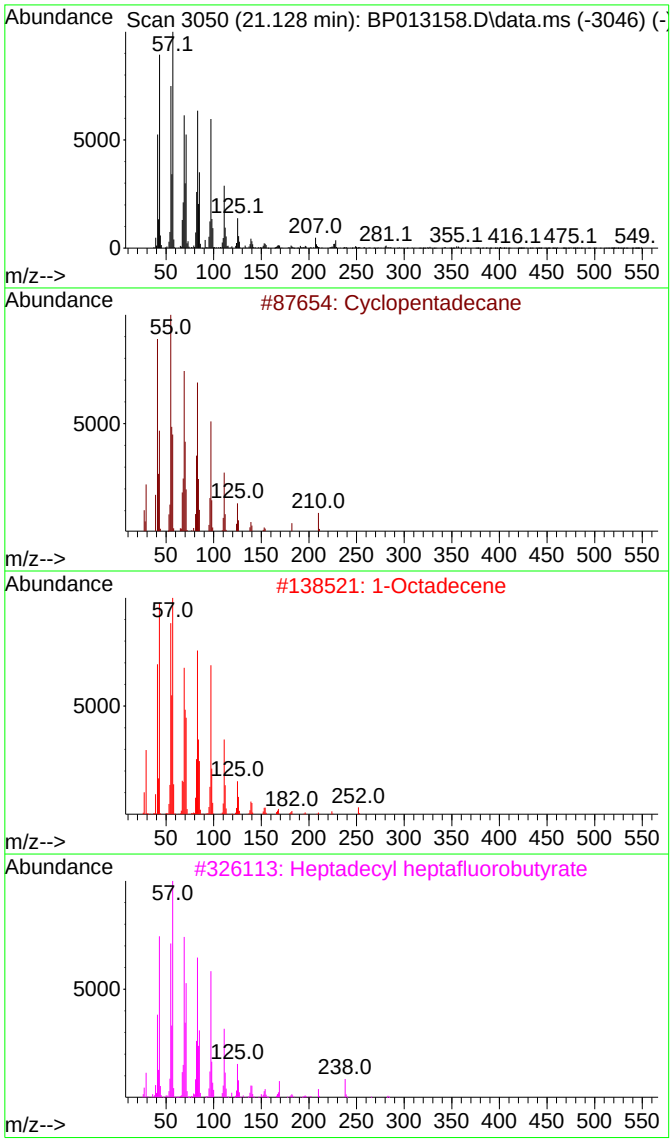
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

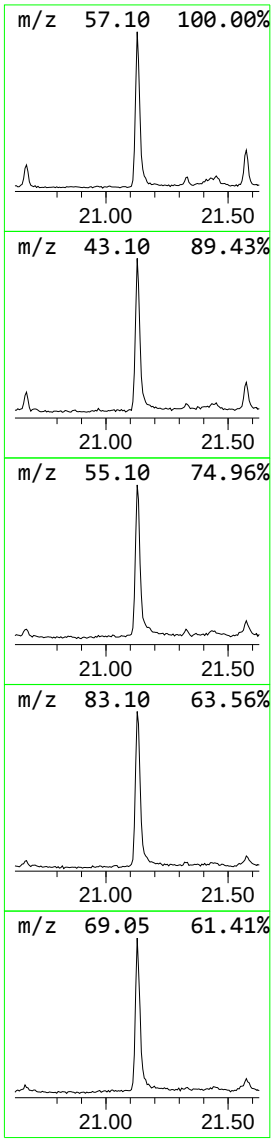
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic21.128 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.128	5.12 ng/ul	1500200	Chrysene-d12	21.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	97
2	1-Octadecene	252	C18H36	000112-88-9	96
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
5	1-Heneicosanol	312	C21H44O	015594-90-8	93





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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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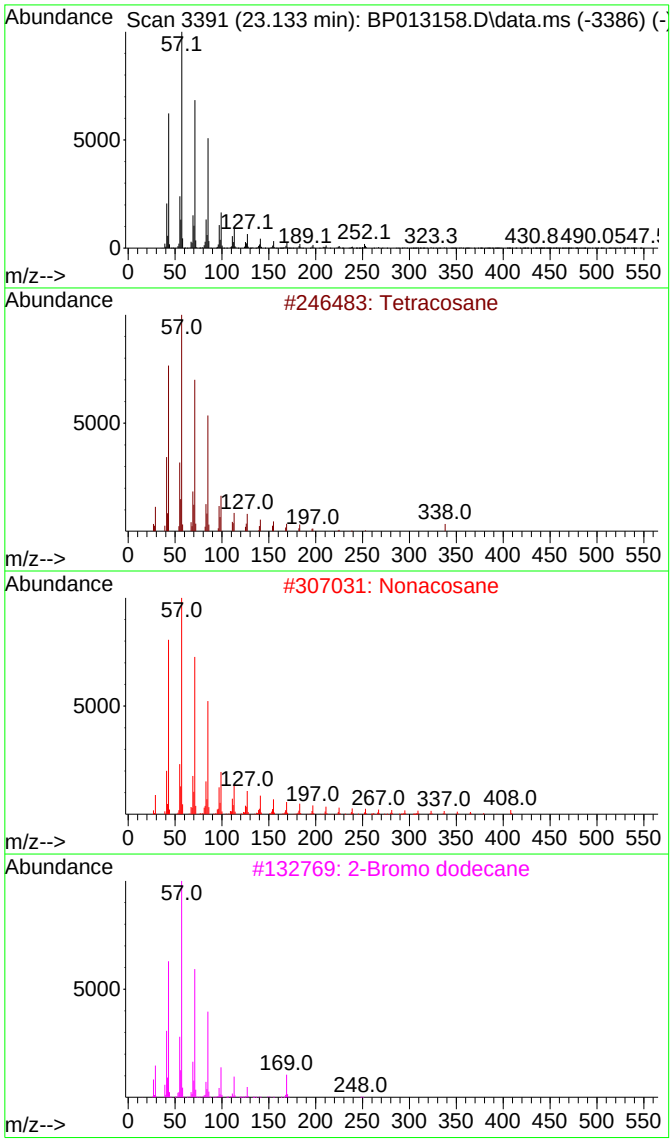
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

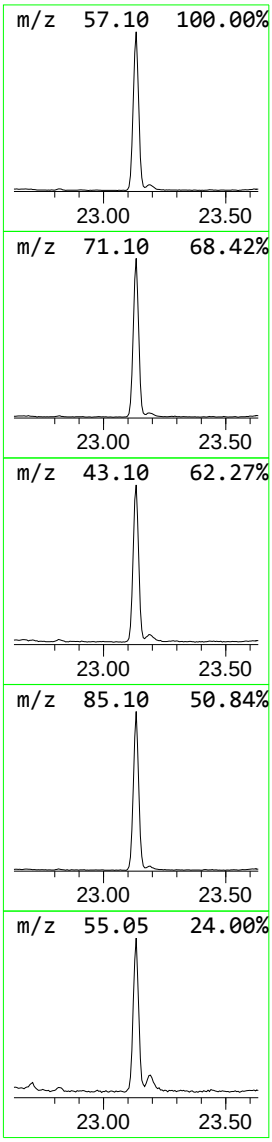
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.133	13.02 ng/ul	4075690	Perylene-d12	23.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	95
2	Nonacosane	408	C29H60	000630-03-5	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Octacosane	394	C28H58	000630-02-4	91





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Data File : BP013158.D
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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

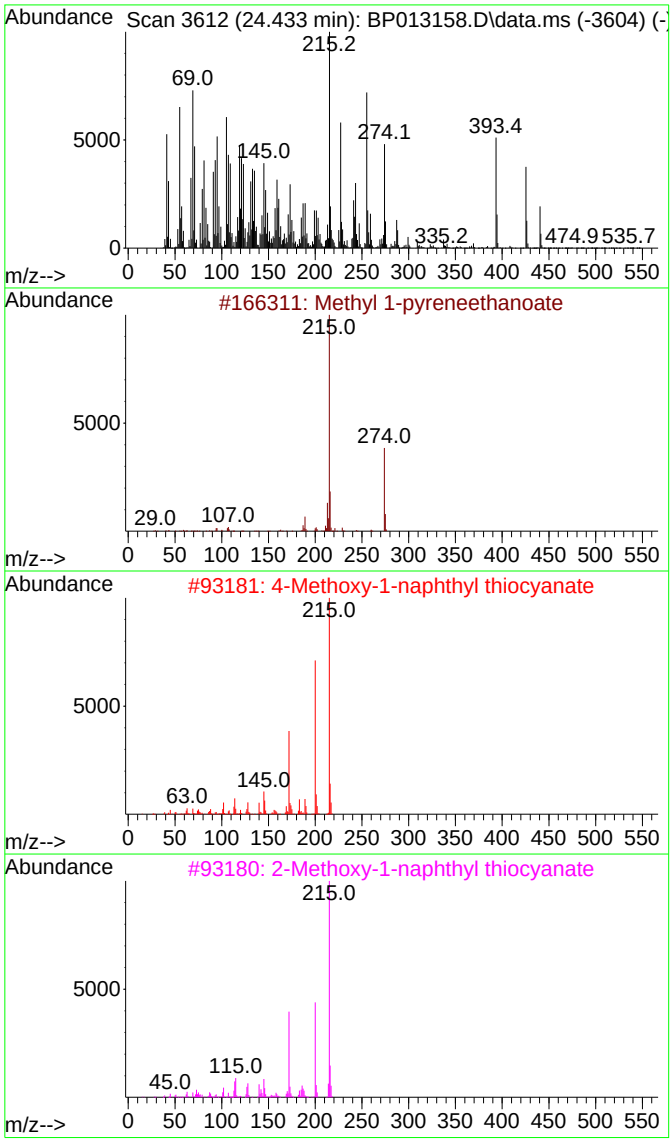
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

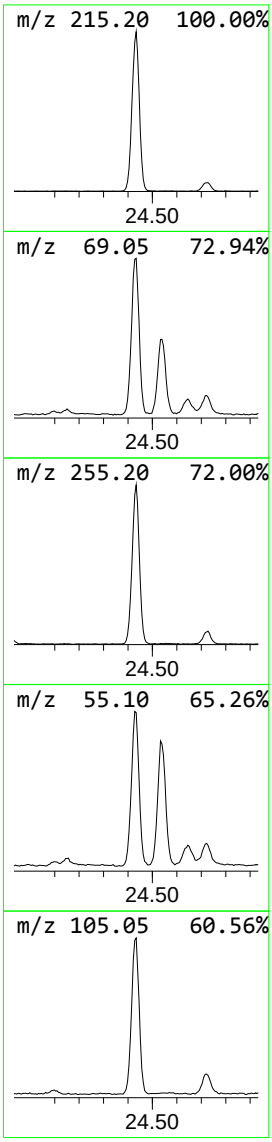
****

Peak Number 8 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.433	38.80 ng/ul	12144300	Perylene-d12	23.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	42
2	4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	18
3	2-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	092696-42-9	18
4	2(3H)-Benzofuranone, 3-[3-(dimet...	215	C13H13NO2	056771-83-6	18
5	4-(1H-Indol-3-yl)-1,3-thiazol-2-...	215	C11H9N3S	022258-56-6	18





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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

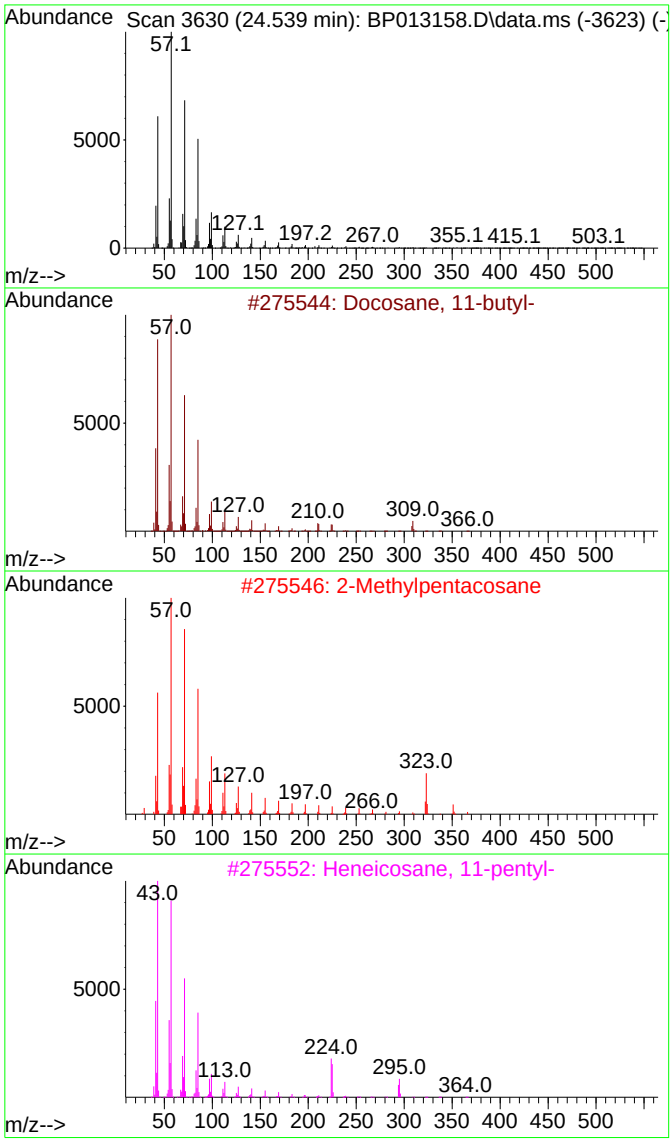
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

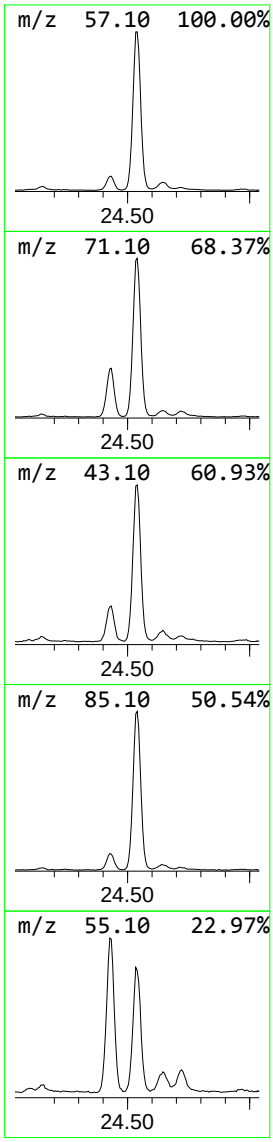
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.539	18.49 ng/ul	5788090	Perylene-d12	23.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2	2-Methylpentacosane	366	C26H54	000629-87-8	95
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
5	Triacontane	422	C30H62	000638-68-6	91





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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

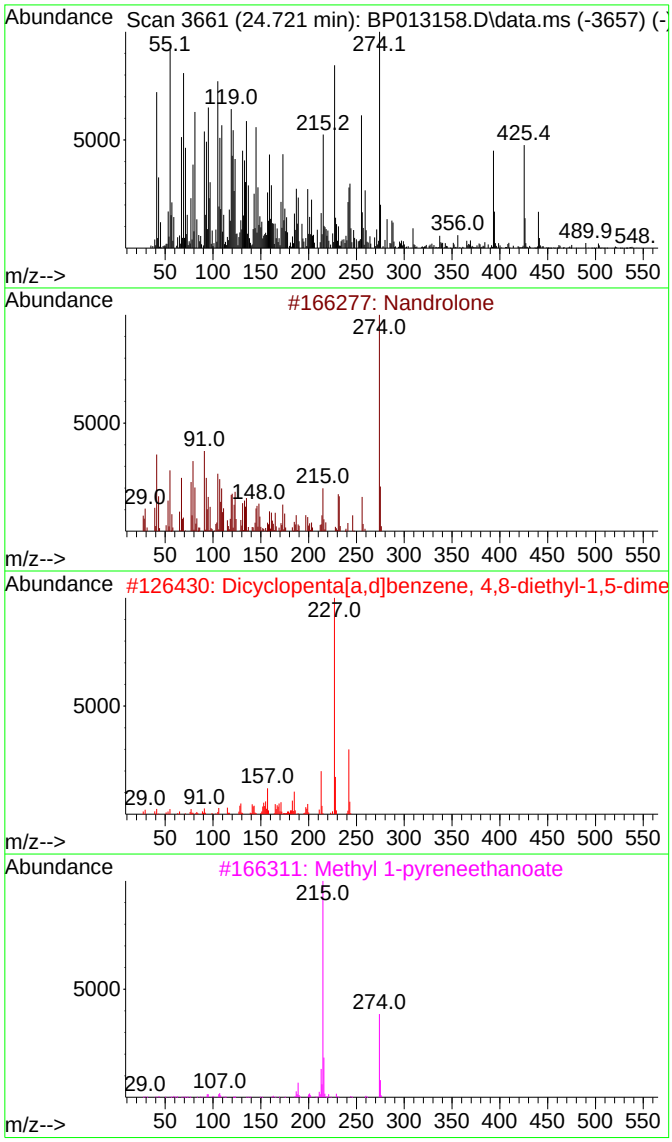
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

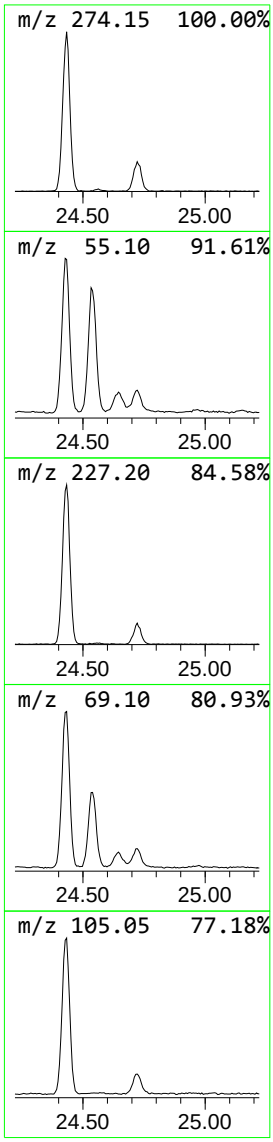
****

Peak Number 10 Nandrolone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.721	4.92 ng/ul	1541430	Perylene-d12	23.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nandrolone	274	C18H26O2	000434-22-0	51
2	Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	44
3	Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	38
4	Androstan-17-one	274	C19H30O	036378-49-1	35
5	Androst-5-en-17-ol	274	C19H30O	1000193-02-3	35





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Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

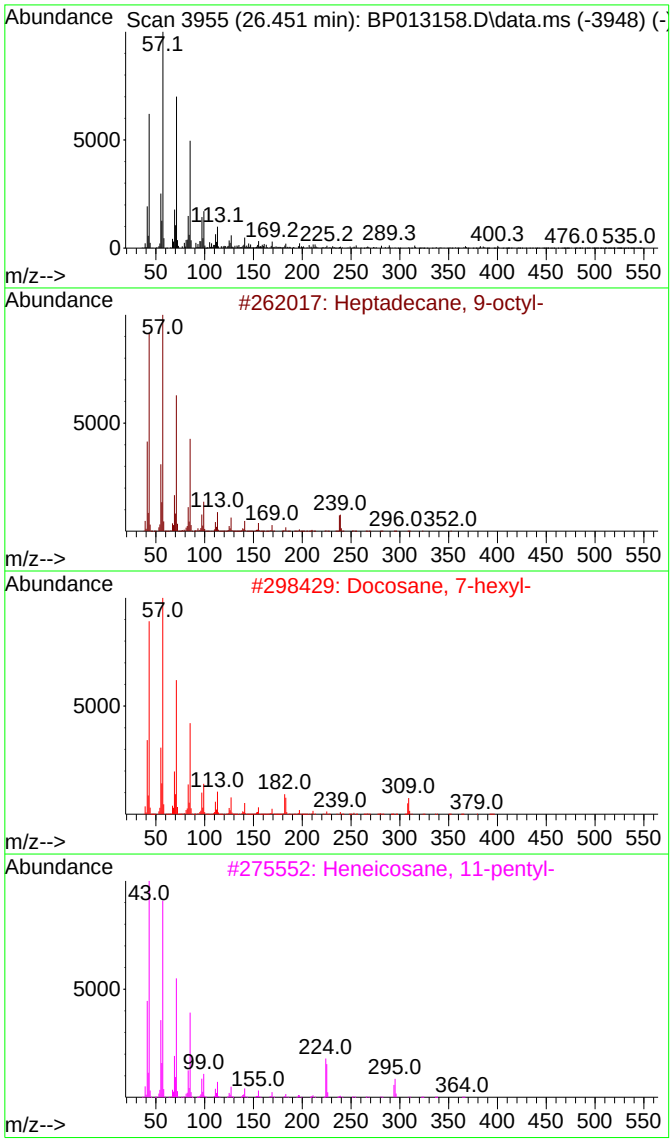
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

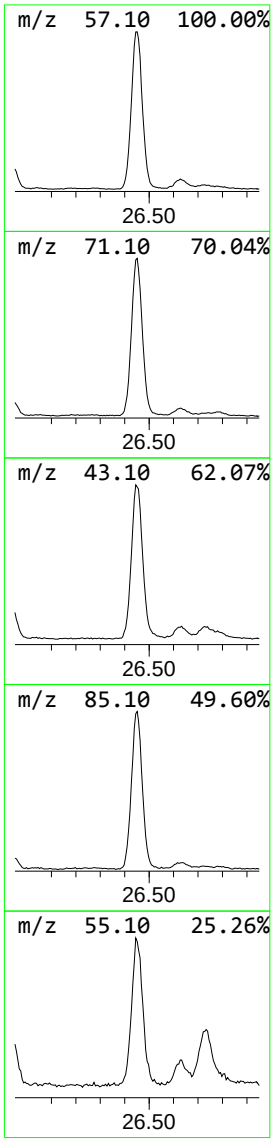
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.451	16.10 ng/ul	5039130	Perylene-d12	23.916

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
4	Docosane, 9-butyl-	366	C26H54	055282-14-9	93
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93





Tentatively Identified Compound (LSC) summary

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Data File : BP013158.D
Acq On : 20 Dec 2022 10:14
Operator : CG/JU
Sample : N6006-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXZ2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-...	5.040	3.0	ng/ul	368614	1	7.952	2433680	20.0
Benzeneacetonit...	12.457	4.6	ng/ul	788683	2	10.769	3444550	20.0
Benzophenone	15.957	17.2	ng/ul	3723000	3	14.593	4318190	20.0
n-Hexadecanoic ...	18.222	4.4	ng/ul	1071340	4	17.351	4907640	20.0
4-(3-Fluoro-8-n...	19.834	4.4	ng/ul	1285180	5	21.439	5859240	20.0
(DEL) Alkane: C...	21.128	5.1	ng/ul	1500200	5	21.439	5859240	20.0
(DEL) Alkane: S...	23.133	13.0	ng/ul	4075690	6	23.916	6259660	20.0
unknown-01	24.433	38.8	ng/ul	12144300	6	23.916	6259660	20.0
(DEL) Alkane: S...	24.539	18.5	ng/ul	5788090	6	23.916	6259660	20.0
Nandrolone	24.721	4.9	ng/ul	1541430	6	23.916	6259660	20.0
(DEL) Alkane: S...	26.451	16.1	ng/ul	5039130	6	23.916	6259660	20.0