

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091224\
 Data File : BN033884.D
 Acq On : 12 Sep 2024 16:50
 Operator : JU/RC
 Sample : P3878-19DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DDCC2DL

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_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN033884.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	72	76	81	rVB	18543	24266	0.76%	0.084%
2	3.505	84	88	96	rBV	42847	65380	2.05%	0.227%
3	4.111	187	191	197	rVV	17587	26294	0.83%	0.091%
4	4.258	208	216	223	rBV	8188	20275	0.64%	0.070%
5	4.370	231	235	240	rVV2	12318	20703	0.65%	0.072%
6	4.687	283	289	300	rBV	446252	589748	18.53%	2.047%
7	5.558	431	437	444	rBV3	14310	23145	0.73%	0.080%
8	5.864	483	489	499	rVV	630051	928754	29.18%	3.223%
9	6.511	595	599	605	rVB3	19958	30717	0.96%	0.107%
10	6.569	605	609	612	rBV2	16162	21947	0.69%	0.076%
11	6.681	618	628	635	rBV2	166712	282002	8.86%	0.979%
12	6.840	650	655	661	rVV	98171	151440	4.76%	0.526%
13	6.940	667	672	679	rVV	52396	82726	2.60%	0.287%
14	7.028	682	687	694	rVV	129482	207443	6.52%	0.720%
15	7.134	699	705	712	rVB2	96973	149140	4.69%	0.518%
16	7.493	758	766	773	rVV	647417	1005489	31.59%	3.490%
17	7.569	773	779	783	rVV4	16145	27399	0.86%	0.095%
18	7.763	809	812	823	rVB8	16853	42677	1.34%	0.148%
19	7.916	830	838	844	rBV	285045	437917	13.76%	1.520%
20	7.975	844	848	853	rVV	55777	76423	2.40%	0.265%
21	8.028	853	857	863	rVB3	30163	49752	1.56%	0.173%
22	8.087	863	867	870	rBV3	20091	31799	1.00%	0.110%
23	8.122	870	873	877	rVV4	18843	33572	1.05%	0.117%
24	8.158	877	879	881	rVV2	22379	25513	0.80%	0.089%
25	8.199	881	886	893	rVB	181322	278456	8.75%	0.966%
26	8.263	893	897	899	rBV	25686	39613	1.24%	0.137%
27	8.293	899	902	909	rVB2	39737	68825	2.16%	0.239%
28	8.369	909	915	922	rVB3	13751	28100	0.88%	0.098%
29	8.434	922	926	930	rBV2	10741	17018	0.53%	0.059%
30	8.487	930	935	945	rVB2	50148	75953	2.39%	0.264%
31	8.634	945	960	966	rBV2	113385	250700	7.88%	0.870%
32	8.722	970	975	980	rVB	87121	120030	3.77%	0.417%
33	8.910	1000	1007	1011	rBV4	13252	27465	0.86%	0.095%
34	8.975	1011	1018	1024	rVB9	9211	21387	0.67%	0.074%
35	9.063	1024	1033	1039	rBV3	43962	85666	2.69%	0.297%
36	9.205	1052	1057	1064	rVB2	40335	73924	2.32%	0.257%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

37	9.346	1075	1081	1090	rBV	114574	192073	6.03%	0.667%
38	9.516	1106	1110	1115	rBV4	22046	32173	1.01%	0.112%
39	9.569	1115	1119	1125	rVB3	16081	29377	0.92%	0.102%
40	9.775	1149	1154	1159	rBV	79698	122906	3.86%	0.427%
41	9.881	1165	1172	1180	rVB	171854	301473	9.47%	1.046%
42	10.093	1204	1208	1212	rVB2	14399	21349	0.67%	0.074%
43	10.157	1212	1219	1228	rVB7	13320	37285	1.17%	0.129%
44	10.252	1228	1235	1246	rBV2	957663	1889495	59.36%	6.558%
45	10.393	1252	1259	1266	rBV2	261635	447420	14.06%	1.553%
46	10.475	1269	1273	1279	rVV2	41250	66997	2.10%	0.233%
47	10.540	1279	1284	1285	rVV2	19391	26330	0.83%	0.091%
48	10.569	1285	1289	1297	rVB	41609	67598	2.12%	0.235%
49	10.646	1297	1302	1311	rVB2	22615	53546	1.68%	0.186%
50	10.781	1321	1325	1329	rBV2	26783	39520	1.24%	0.137%
51	11.157	1384	1389	1393	rBV3	26743	49503	1.56%	0.172%
52	11.216	1396	1399	1403	rVV	17996	28642	0.90%	0.099%
53	11.304	1409	1414	1420	rVB4	13801	24275	0.76%	0.084%
54	11.504	1441	1448	1457	rBV6	22483	47982	1.51%	0.167%
55	11.687	1474	1479	1489	rVB4	35359	63743	2.00%	0.221%
56	11.799	1489	1498	1506	rBV3	32240	82319	2.59%	0.286%
57	11.928	1515	1520	1524	rVB4	15525	24055	0.76%	0.083%
58	12.040	1528	1539	1544	rBV2	41190	86751	2.73%	0.301%
59	12.122	1546	1553	1562	rVB5	57609	111007	3.49%	0.385%
60	12.304	1579	1584	1588	rBV4	19605	34779	1.09%	0.121%
61	12.551	1618	1626	1635	rBV	130368	209016	6.57%	0.725%
62	12.657	1635	1644	1647	rVV4	19531	39907	1.25%	0.139%
63	12.704	1647	1652	1659	rVV	102376	157408	4.94%	0.546%
64	12.769	1659	1663	1673	rVB7	13385	31923	1.00%	0.111%
65	12.975	1693	1698	1703	rVV6	12371	23597	0.74%	0.082%
66	13.134	1721	1725	1728	rVV5	14144	25006	0.79%	0.087%
67	13.193	1731	1735	1740	rVB2	32811	53321	1.68%	0.185%
68	13.328	1746	1758	1762	rBV7	26647	60625	1.90%	0.210%
69	13.557	1792	1797	1803	rBV	346225	450860	14.16%	1.565%
70	13.822	1836	1842	1850	rBV2	345187	511947	16.08%	1.777%
71	13.898	1850	1855	1860	rVB2	9396	18894	0.59%	0.066%
72	13.951	1860	1864	1868	rBV5	11606	18267	0.57%	0.063%
73	14.069	1877	1884	1887	rBV4	14902	23391	0.73%	0.081%
74	14.140	1889	1896	1904	rVB	1220018	1798736	56.51%	6.243%
75	14.357	1926	1933	1946	rBV	132575	251529	7.90%	0.873%
76	14.469	1946	1952	1957	rVV8	9865	25550	0.80%	0.089%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

77	14.775	1998	2004	2009	rVV	217572	313713	9.86%	1.089%
78	15.028	2043	2047	2052	rVB2	30255	39362	1.24%	0.137%
79	15.140	2060	2066	2076	rVB	462414	680868	21.39%	2.363%
80	15.263	2078	2087	2090	rBV	125611	194524	6.11%	0.675%
81	15.292	2090	2092	2104	rVB5	34925	61483	1.93%	0.213%
82	15.516	2122	2130	2136	rVB	48608	74542	2.34%	0.259%
83	15.892	2191	2194	2198	rBV2	16844	24946	0.78%	0.087%
84	15.969	2203	2207	2212	rVB2	19448	27995	0.88%	0.097%
85	16.545	2298	2305	2318	rBV	1704615	2338295	73.46%	8.116%
86	16.645	2320	2322	2325	rVV4	13032	20586	0.65%	0.071%
87	16.687	2326	2329	2333	rVV4	22421	36428	1.14%	0.126%
88	16.892	2354	2364	2370	rBV	1601432	2212273	69.50%	7.678%
89	16.992	2374	2381	2387	rBV	539961	772759	24.28%	2.682%
90	17.428	2452	2455	2459	rBV2	15315	18870	0.59%	0.065%
91	17.522	2467	2471	2475	rVB2	35971	43271	1.36%	0.150%
92	17.822	2517	2522	2531	rBV	323831	473900	14.89%	1.645%
93	18.122	2569	2573	2579	rVB3	14269	20644	0.65%	0.072%
94	18.763	2679	2682	2690	rVB7	16076	25163	0.79%	0.087%
95	19.116	2738	2742	2752	rBV	114990	172257	5.41%	0.598%
96	19.298	2767	2773	2780	rVV	735156	948081	29.78%	3.291%
97	20.857	3034	3038	3045	rBV4	131912	210473	6.61%	0.730%
98	21.127	3079	3084	3097	rBV	1967296	2795698	87.83%	9.703%
99	23.716	3517	3524	3536	rVB	485397	1100739	34.58%	3.820%
100	23.910	3548	3557	3565	rBV	1356895	3183183	100.00%	11.048%

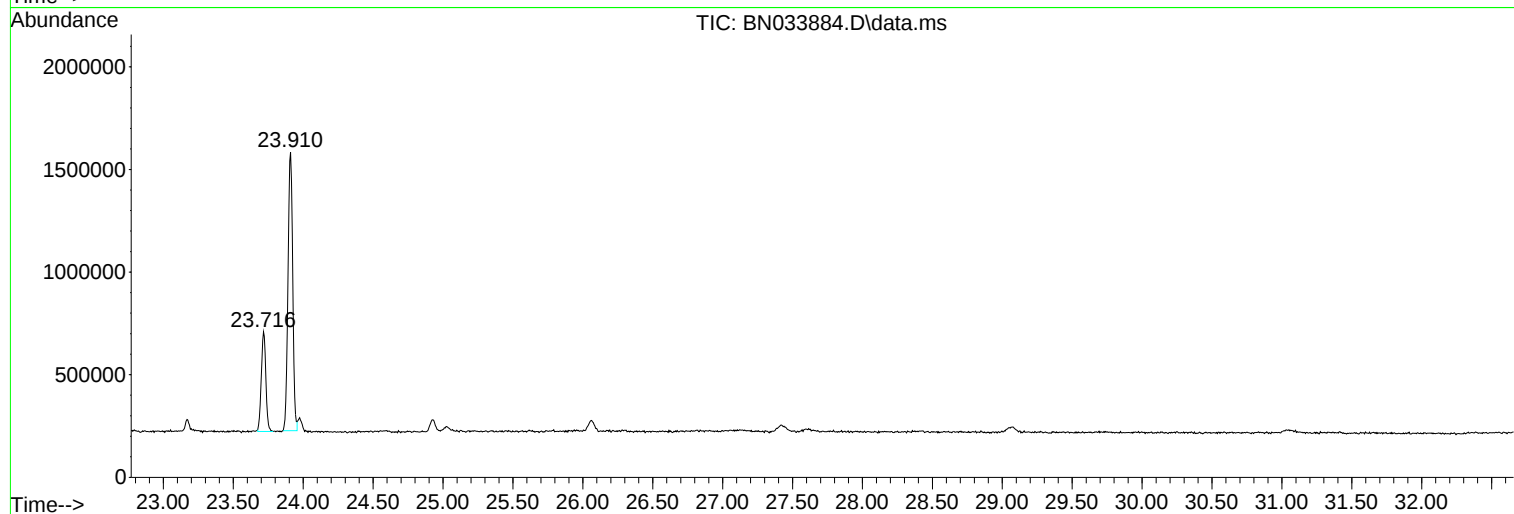
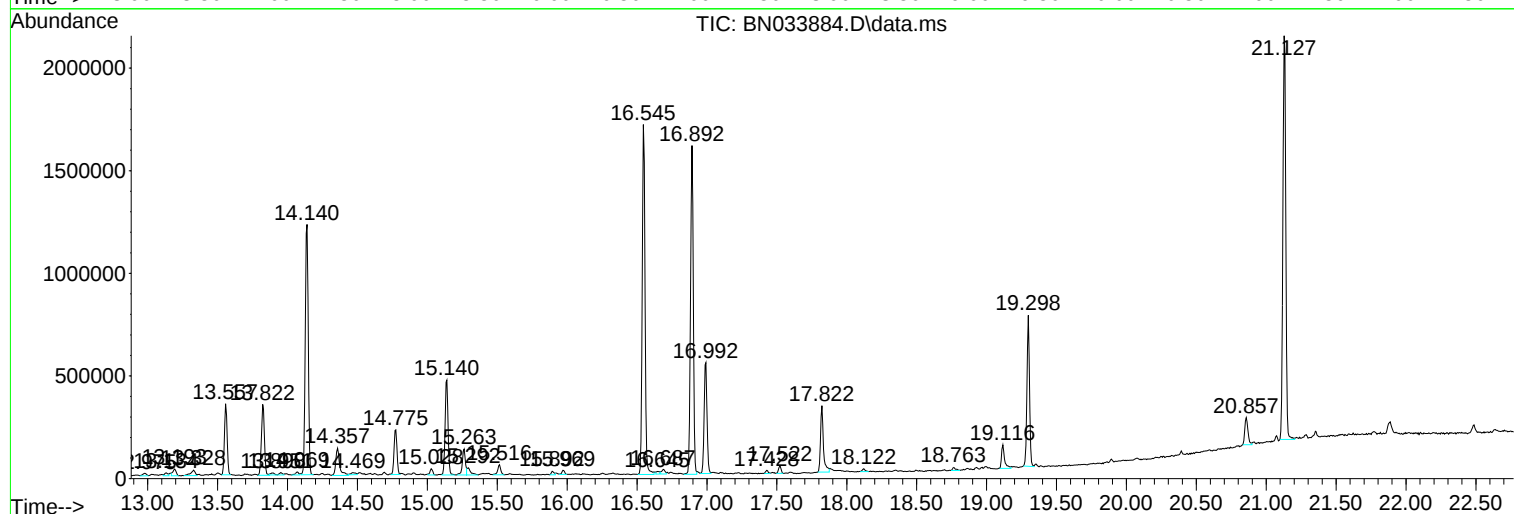
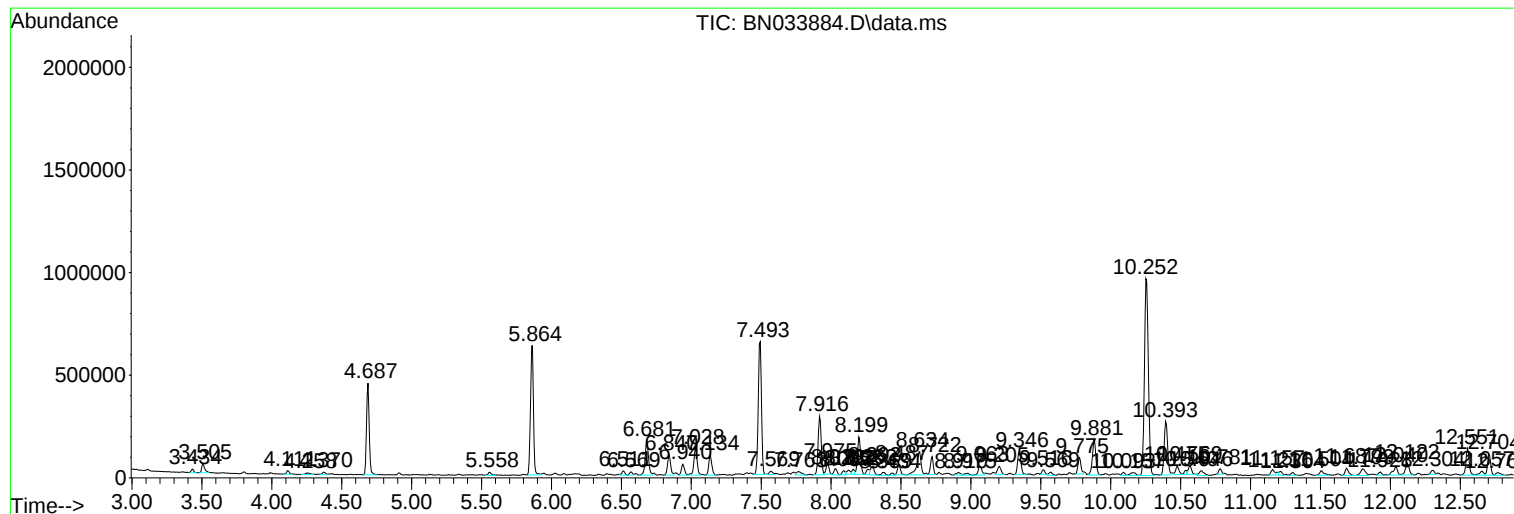
Sum of corrected areas: 28812286

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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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



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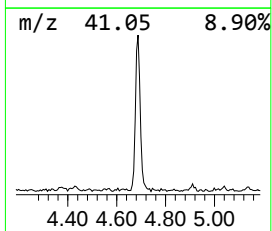
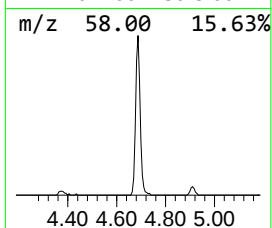
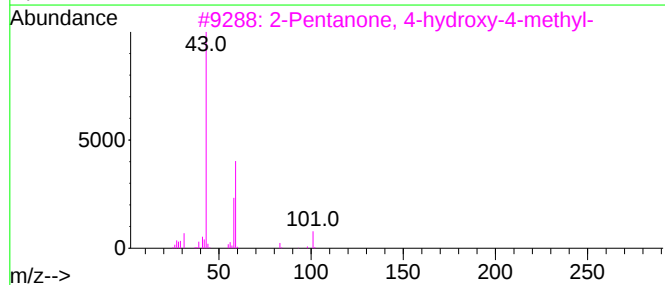
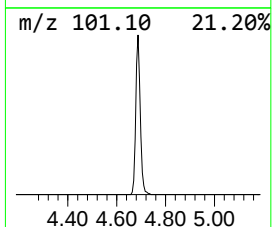
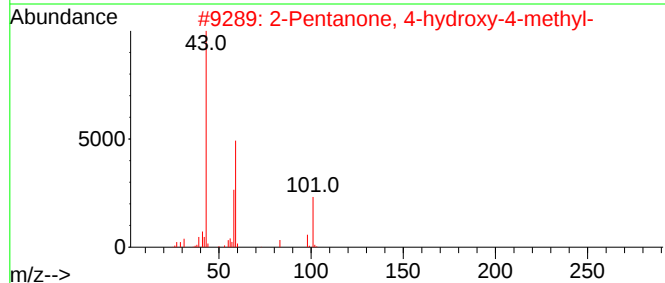
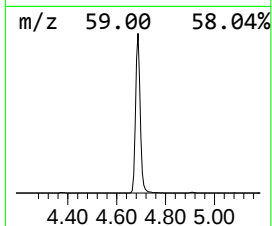
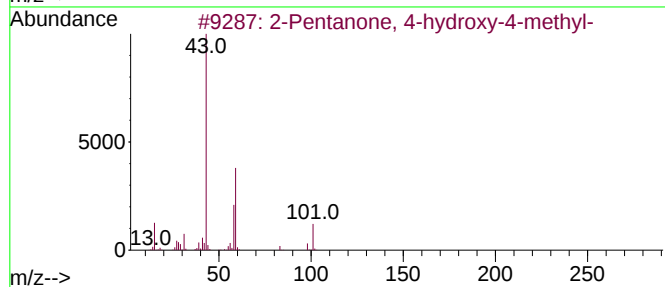
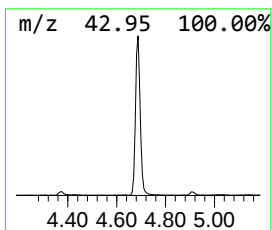
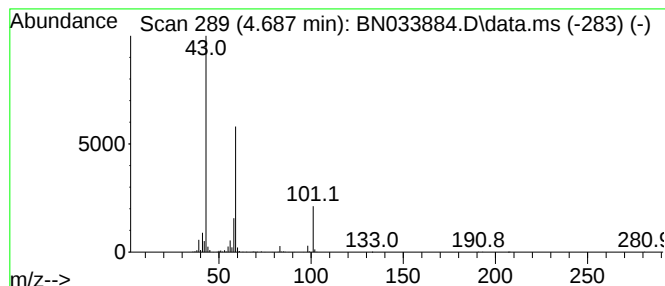
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.687	11.73 ng/ul	589748	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		



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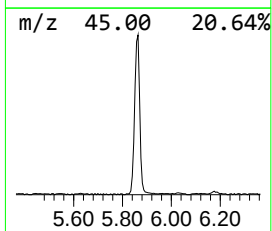
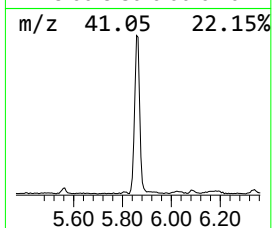
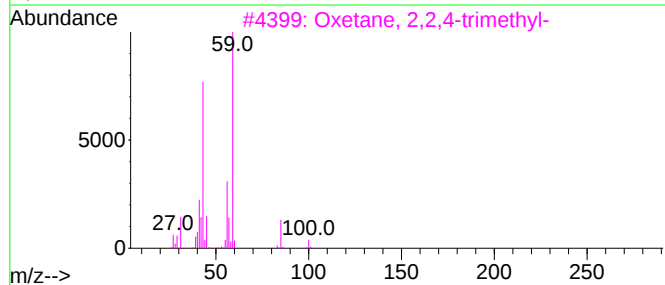
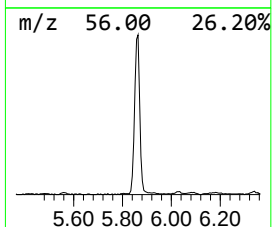
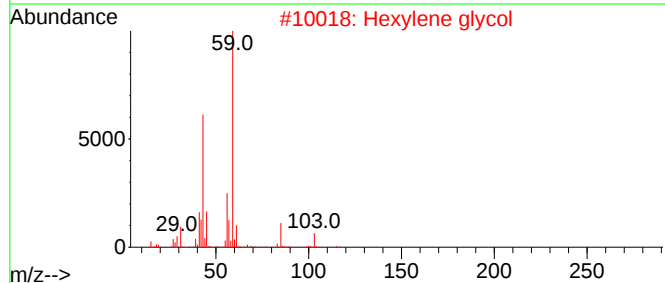
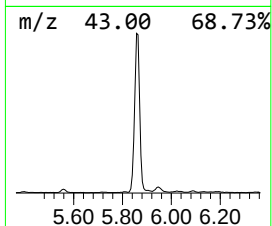
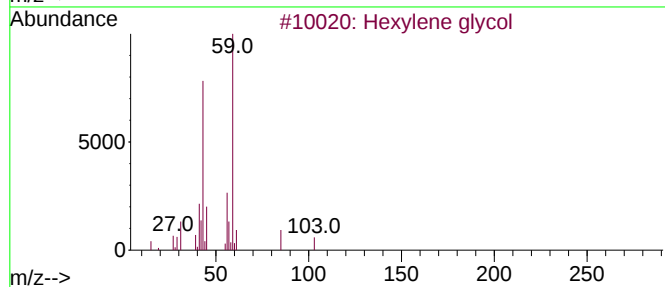
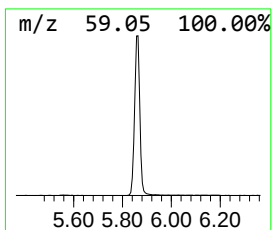
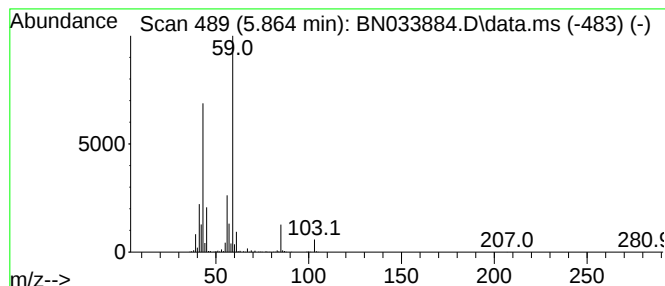
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.864	18.47 ng/ul	928754	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	90
2			Hexylene glycol	118	C6H14O2	000107-41-5	90
3			Oxetane, 2,2,4-trimethyl-	100	C6H12O	023120-44-7	72
4			(R)-(-)-2-Methyl-2,4-pentanediol	118	C6H14O2	099210-90-9	59
5			Hexylene glycol	118	C6H14O2	000107-41-5	50



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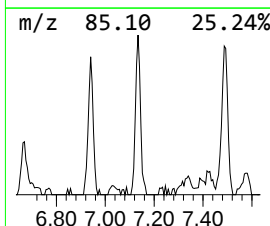
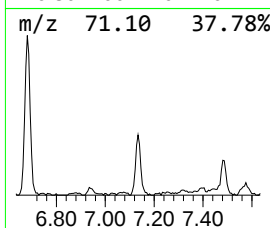
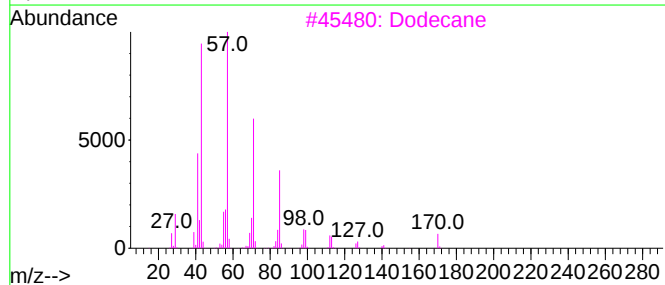
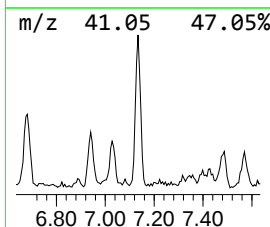
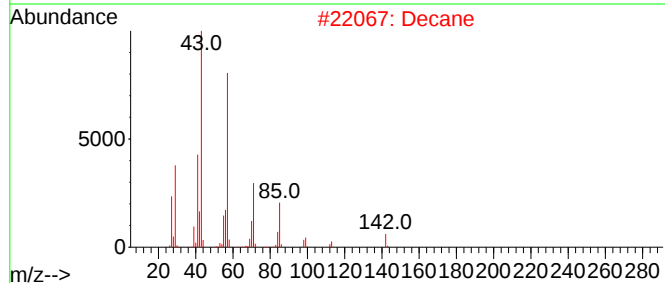
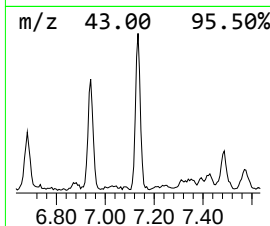
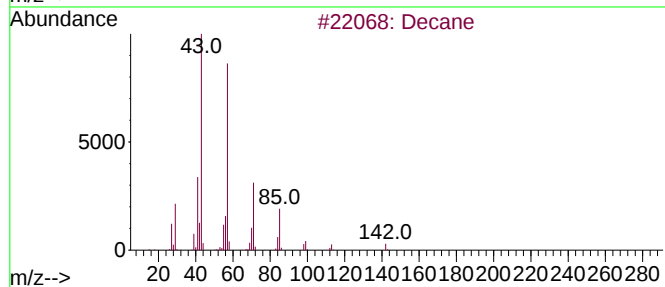
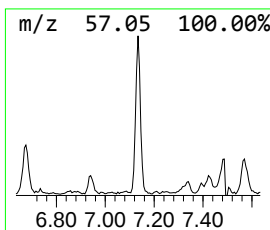
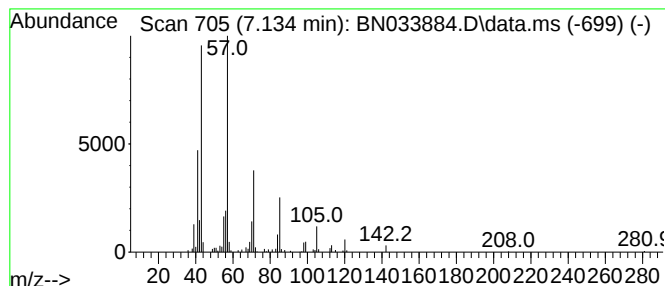
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.134	2.97 ng/ul	149140	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	94
2			Decane	142	C10H22	000124-18-5	80
3			Dodecane	170	C12H26	000112-40-3	80
4			Tridecane	184	C13H28	000629-50-5	72
5			Undecane	156	C11H24	001120-21-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091224\
Data File : BN033884.D
Acq On : 12 Sep 2024 16:50
Operator : JU/RC
Sample : P3878-19DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC2DL

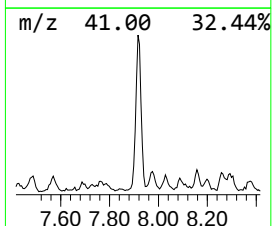
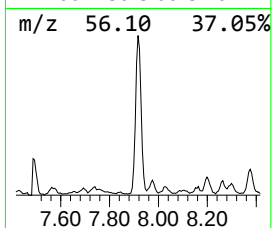
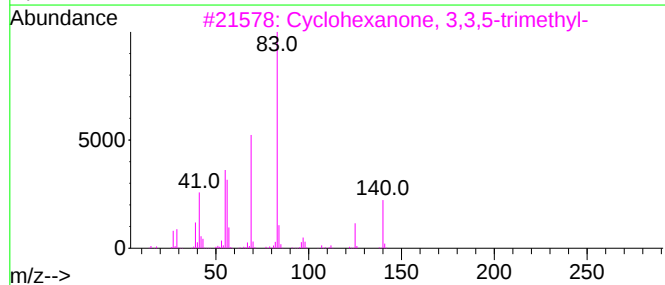
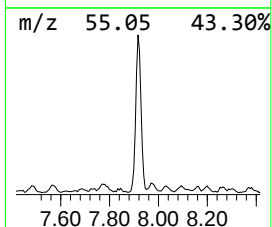
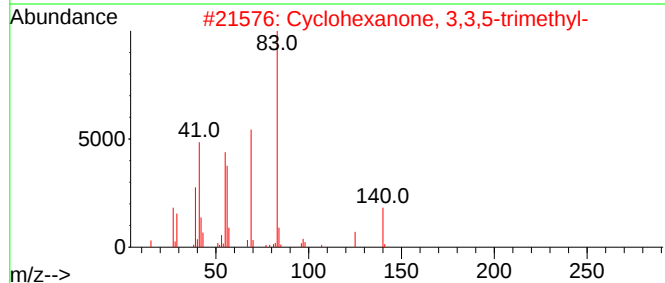
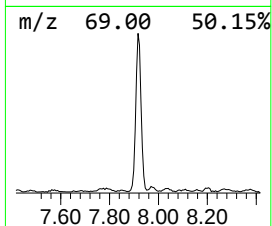
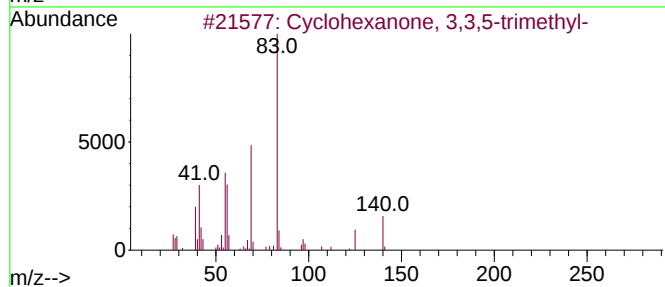
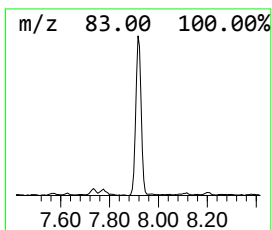
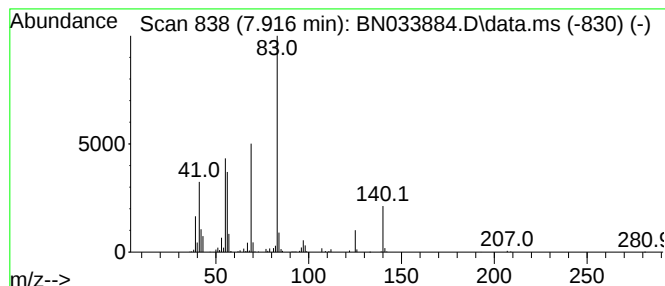
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexanone, 3,3,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.916	8.71 ng/ul	437917	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5			1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091224\
Data File : BN033884.D
Acq On : 12 Sep 2024 16:50
Operator : JU/RC
Sample : P3878-19DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC2DL

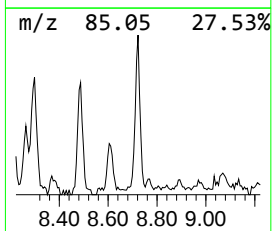
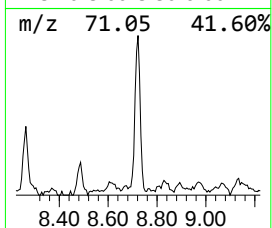
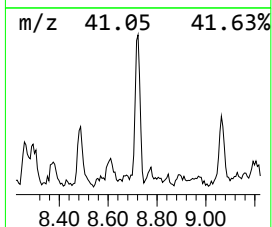
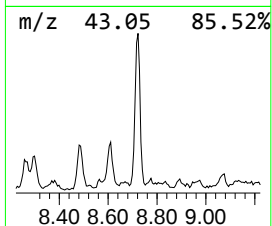
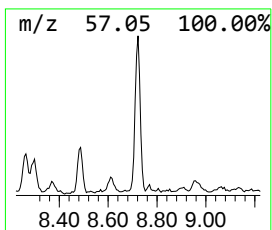
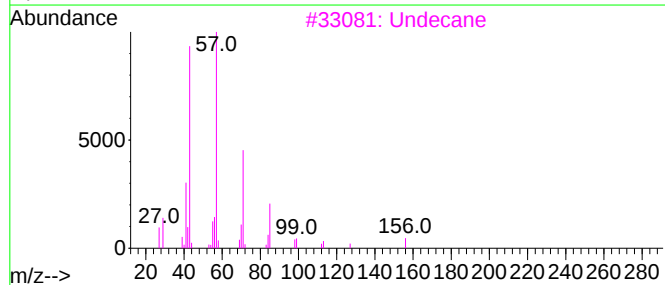
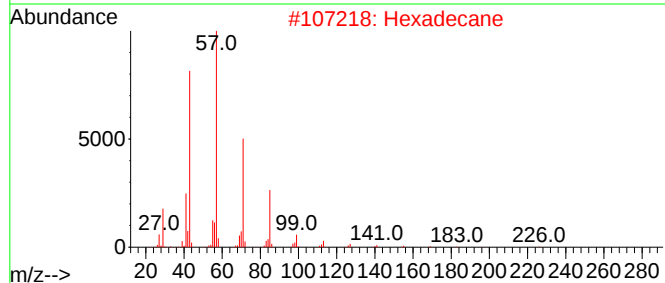
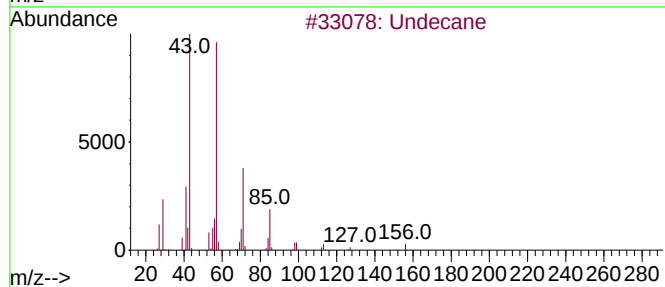
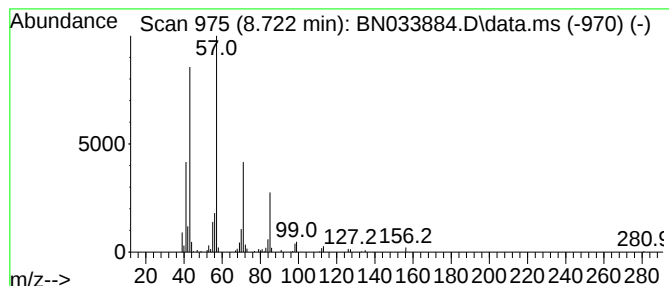
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.722	2.39 ng/ul	120030	1,4-Dichlorobenzene-d4	7.493

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	94
2			Hexadecane	226	C16H34	000544-76-3	90
3			Undecane	156	C11H24	001120-21-4	83
4			Tridecane	184	C13H28	000629-50-5	78
5			Pentadecane	212	C15H32	000629-62-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091224\
Data File : BN033884.D
Acq On : 12 Sep 2024 16:50
Operator : JU/RC
Sample : P3878-19DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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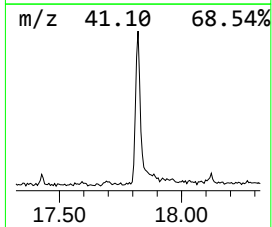
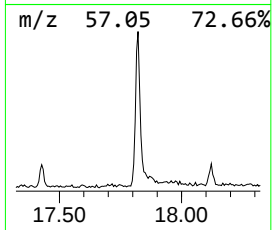
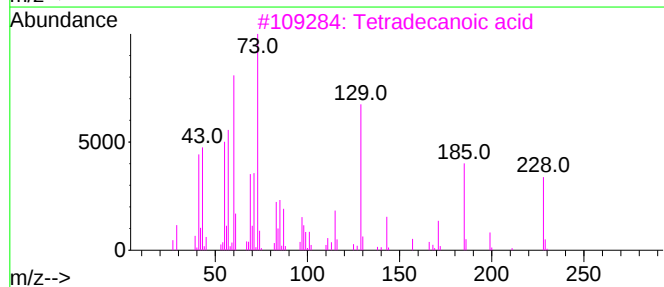
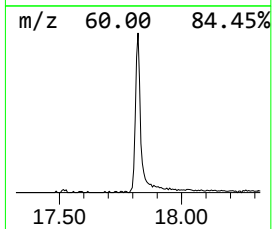
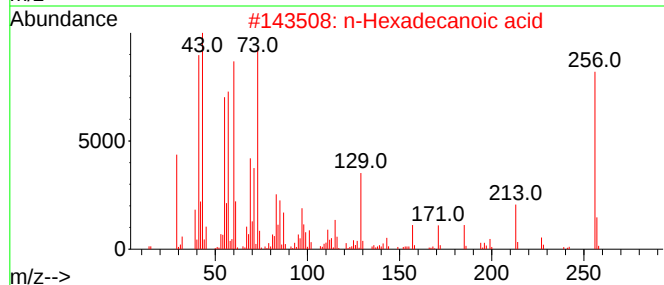
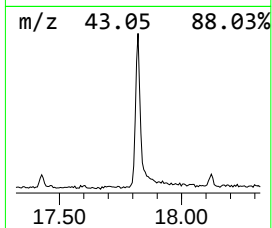
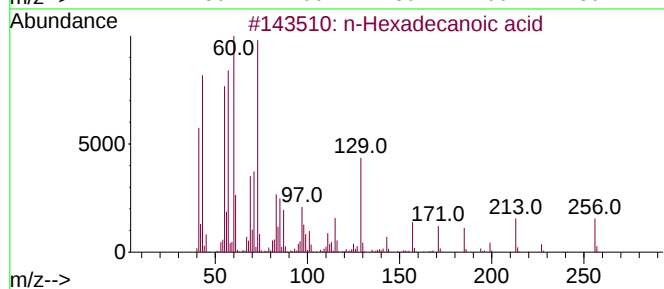
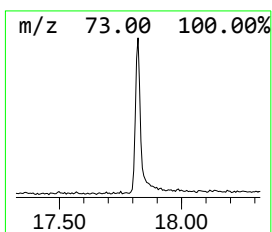
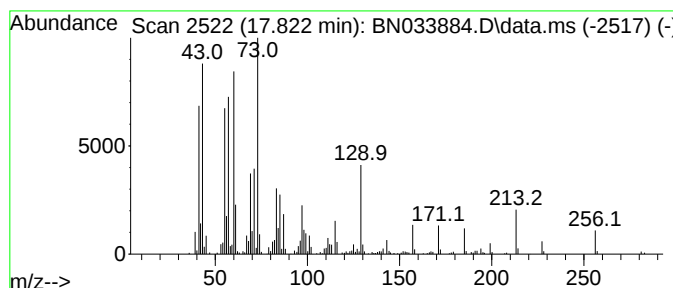
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.822	4.28 ng/ul	473900	Phenanthrene-d10	16.892

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	Tetradecanoic acid			228	C14H28O2	000544-63-8	91
4	Pentadecanoic acid			242	C15H30O2	001002-84-2	90
5	Tetradecanoic acid			228	C14H28O2	000544-63-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091224\
Data File : BN033884.D
Acq On : 12 Sep 2024 16:50
Operator : JU/RC
Sample : P3878-19DL 5X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.687	11.7	ng/ul	589748	1	7.493	1005490	20.0
Hexylene glycol	5.864	18.5	ng/ul	928754	1	7.493	1005490	20.0
(DEL) Alkane: S...	7.134	3.0	ng/ul	149140	1	7.493	1005490	20.0
Cyclohexanone, ...	7.916	8.7	ng/ul	437917	1	7.493	1005490	20.0
(DEL) Alkane: S...	8.722	2.4	ng/ul	120030	1	7.493	1005490	20.0
n-Hexadecanoic ...	17.822	4.3	ng/ul	473900	4	16.892	2212270	20.0