LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

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-BP062322.M
Title : SVOA CALIBRATION

Signal : TIC: BP010809.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.246	23	27	38	rVB2	360104	477044	0.21%	0.051%
2	3.646	88	95	103	rBV3	225424	938284	0.41%	0.100%
3	6.234	529	535	546	rVV	976752	1713098	0.74%	0.183%
4	6.928	643	653	665	rVB	13358105	22902239	9.94%	2.442%
5	7.063	670	676	686	rVV	1078697	1622231	0.70%	0.173%
6	7.252	703	708	716	rVB	1369627	2112286	0.92%	0.225%
7	7.722	780	788	794	rBV	647917	1053095	0.46%	0.112%
8	7.805	794	802	815	rBV	1857442	5222939	2.27%	0.557%
9	7.946	821	826	833	rVB	392086	661702	0.29%	0.071%
10	8.399	892	903	905	rBV2	403784	967812	0.42%	0.103%
11	8.434	905	909	914	rVB	980249	1494745	0.65%	0.159%
12	8.499	914	920	925	rBV	1099948	1982899	0.86%	0.211%
13	8.881	978	985	991	rVV	1027185	1785165	0.77%	0.190%
14	9.099	1012	1022	1027	rBV	395374	995214	0.43%	0.106%
15	9.222	1036	1043	1047	rBV6	221133	534133	0.23%	0.057%
16	9.334	1057	1062	1070	rBV	446588	797228	0.35%	0.085%
17	9.416	1070	1076	1084	rBV4	370383	1065630	0.46%	0.114%
18	9.499	1084	1090	1097	rVB	966702	1671168	0.72%	0.178%
19	9.569	1097	1102	1104	rBV2	599603	991758	0.43%	0.106%
20	9.599	1104	1107	1113	rVB	955523	1509931	0.66%	0.161%
21	9.704	1118	1125	1128	rBV	4356822	7986349	3.46%	0.852%
22	9.734	1128	1130	1137	rVB2	4427674	6451399	2.80%	0.688%
23	9.916	1155	1161	1164	rBV2	625403	1268399	0.55%	0.135%
24	9.981	1164	1172	1178	rBV2	3550799	10432813	4.53%	1.113%
25	10.034	1178	1181	1187	rVB	2539904	3924549	1.70%	0.419%
26	10.140	1187	1199	1201	rBV2	2554872	5405877	2.35%	0.577%
27	10.228	1210	1214	1218	rBV3	721634	1192520	0.52%	0.127%
28	10.346	1225	1234	1240	rBV7	1285294	4233155	1.84%	0.451%
29	10.404	1241	1244	1246	rVV2	1037207	1563670	0.68%	0.167%
30	10.516	1246	1263	1268	rVV3	46601431	217720886	94.45%	23.220%
31	10.557	1268	1270	1281	rVB3	2441466	4529254	1.96%	0.483%
32	10.710	1286	1296	1301	rBV6	1694213	4291130	1.86%	0.458%
33	10.763	1301	1305	1307	rBV	571145	839464	0.36%	0.090%
34	10.804	1307	1312	1319	rVB3	1909407	3884192	1.69%	0.414%
35	10.910	1319	1330	1336	rBV	29471038	58843126	25.53%	6.276%
36	11.075	1354	1358	1361	rBV	569277	948144	0.41%	0.101%

LSC Area Percent Report

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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

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-BP062322.M
Title : SVOA CALIBRATION

37	11.122	1361	1366	1369	rVB	1065819	1675921	0.73%	0.179%
38	11.257	1383	1389	1394	rBV5	355220	873697	0.38%	0.093%
39	11.322	1394	1400	1405	rBV3	698259	1323364	0.57%	0.141%
40	11.416	1405	1416	1420	rBV10	987550	3416232	1.48%	0.364%
41	11.510	1423	1432	1436	rVB4	1750301	4576911	1.99%	0.488%
42	11.557	1436	1440	1445	rVB3	2990988	5228174	2.27%	0.558%
43	11.604	1445	1448	1452	rVB3	441613	589665	0.26%	0.063%
44	11.675	1452	1460	1464	rBV2	2629221	6702229	2.91%	0.715%
45	11.746	1468	1472	1475	rBV4	1443346	2527381	1.10%	0.270%
46	11.810	1480	1483	1486	rVV	5318435	7387303	3.20%	0.788%
47	11.840	1486	1488	1490	rVV2	1635807	1887047	0.82%	0.201%
48	11.963	1490	1509	1513	rVV3	46755076	230513934	100.00%	24.584%
49	12.022	1515	1519	1521	rVV2	1778175	3104113	1.35%	0.331%
50	12.057	1521	1525	1528	rVV3	2382375	4908019	2.13%	0.523%
51	12.087	1528	1530	1534	rVV2	2440398	2902734	1.26%	0.310%
52	12.157	1534	1542	1546	rVV	1337620	3036112	1.32%	0.324%
53	12.245	1546	1557	1566	rVB6	2241533	7454908	3.23%	0.795%
54	12.510	1594	1602	1607	rBV8	223973	591606	0.26%	0.063%
55	12.734	1632	1640	1643	rBV2	1232630	2326439	1.01%	0.248%
56	12.775	1643	1647	1658	rVB	3166180	4798076	2.08%	0.512%
57	12.945	1669	1676	1683	rBV9	174051	527525	0.23%	0.056%
58	13.257	1718	1729	1733	rBV2	49637582	117881217	51.14%	12.572%
59	13.345	1740	1744	1751	rVB	1393016	1941259	0.84%	0.207%
60	13.604	1781	1788	1792	rVV	26614620	40542658	17.59%	4.324%
61	13.645	1792	1795	1802	rVV	10667844	14031436	6.09%	1.496%
62	13.710	1802	1806	1814	rVV	3140035	4363522	1.89%	0.465%
63	13.810	1814	1823	1827	rVV2	1752015	3118924	1.35%	0.333%
64	13.845	1827	1829	1838	rVB	668328	951127	0.41%	0.101%
65	14.075	1861	1868	1876	rVB	2155615	3514898	1.52%	0.375%
66	14.151	1876	1881	1885	rBV	1358467	1748993	0.76%	0.187%
67	14.245	1893	1897	1901	rBV	1199418	1620588	0.70%	0.173%
68	14.316	1906	1909	1915	rVB2	448456	629583	0.27%	0.067%
69	14.381	1915	1920	1925	rBV2	1113131	1833422	0.80%	0.196%
70	14.434	1925	1929	1935	rVB	1401248	1910115	0.83%	0.204%
71	14.581	1950	1954	1958	rBV2	351786	497582	0.22%	0.053%
72	14.634	1958	1963	1967	rBV	2486268	3243855	1.41%	0.346%
73	14.881	2001	2005	2008	rVB	701578	874570	0.38%	0.093%
74	14.928	2008	2013	2021	rVB3	446349	859679	0.37%	0.092%
75	15.063	2032	2036	2046	rBV	2995379	4099648	1.78%	0.437%
76	15.257	2065	2069	2074	rVB	924249	1143628	0.50%	0.122%

LSC Area Percent Report

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Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

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-BP062322.M
Title : SVOA CALIBRATION

77	15.375	2084	2089	2097	rBV	2623298	4019634	1.74%	0.429%
78	15.492	2105	2109	2115	rBV2	833954	1359359	0.59%	0.145%
79	15.551	2115	2119	2127	rVB	885978	1293368	0.56%	0.138%
80	15.834	2163	2167	2173	rVB	795746	1096878	0.48%	0.117%
81	16.075	2202	2208	2217	rBV	898721	1637255	0.71%	0.175%
82	16.439	2266	2270	2275	rVB	1063208	1319919	0.57%	0.141%
83	16.898	2344	2348	2354	rVB	866720	1326497	0.58%	0.141%
84	17.139	2384	2389	2398	rVB	1230762	1970624	0.85%	0.210%
85	17.239	2401	2406	2418	rVB2	2940714	4269139	1.85%	0.455%
86	17.528	2451	2455	2463	rBV2	376795	710172	0.31%	0.076%
87	18.057	2541	2545	2555	rBV	1000501	1548921	0.67%	0.165%
88	18.263	2576	2580	2587	rBV	1231185	1832688	0.80%	0.195%
89	19.304	2753	2757	2767	rBV2	633809	947517	0.41%	0.101%
90	19.486	2783	2788	2793	rBV	3411388	4473777	1.94%	0.477%
91	19.545	2793	2798	2813	rVB	6599229	8432496	3.66%	0.899%
92	20.098	2888	2892	2902	rBV4	454383	823734	0.36%	0.088%
93	20.363	2934	2937	2941	rBV	667137	724043	0.31%	0.077%
94	20.739	2996	3001	3005	rBV3	525343	797475	0.35%	0.085%
95	20.863	3018	3022	3030	rBV3	935539	1608773	0.70%	0.172%
96	20.986	3039	3043	3049	rVB2	2007967	2432317	1.06%	0.259%
97	21.251	3085	3088	3094	rVB	1616403	1848017	0.80%	0.197%
98	21.533	3132	3136	3141	rVB	623101	801110	0.35%	0.085%
99	23.468	3458	3465	3474	rVB2	2715380	5240158	2.27%	0.559%
100	23.621	3486	3491	3501	rVB2	1004685	1971531	0.86%	0.210%

Sum of corrected areas: 937657024

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
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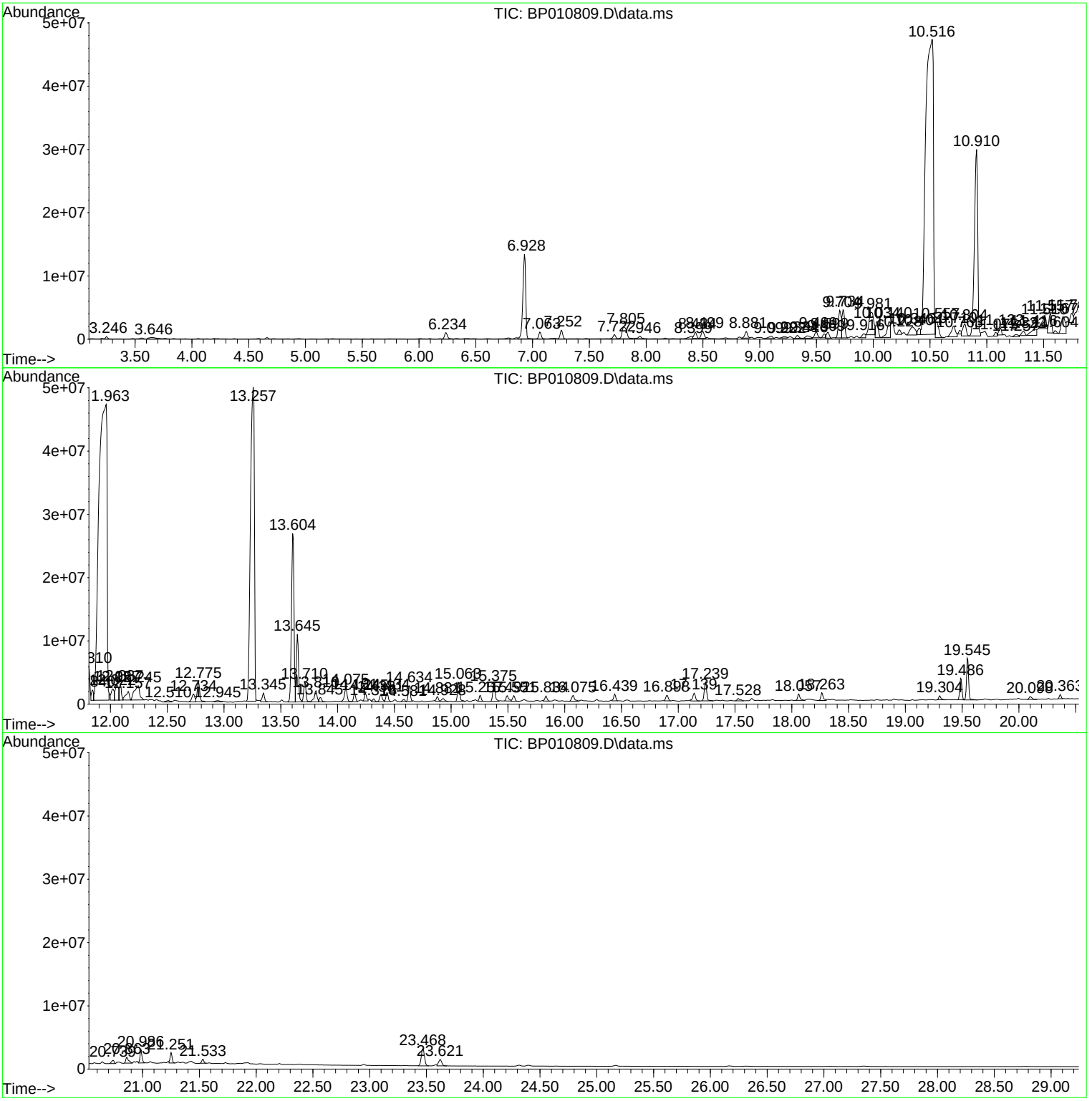
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

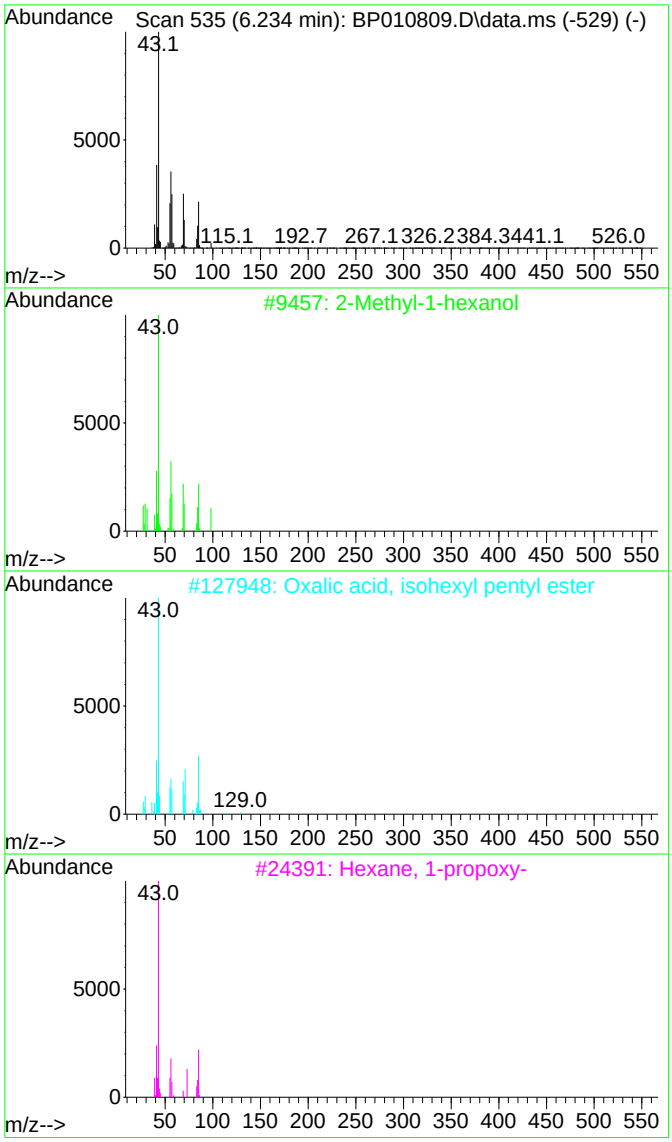
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

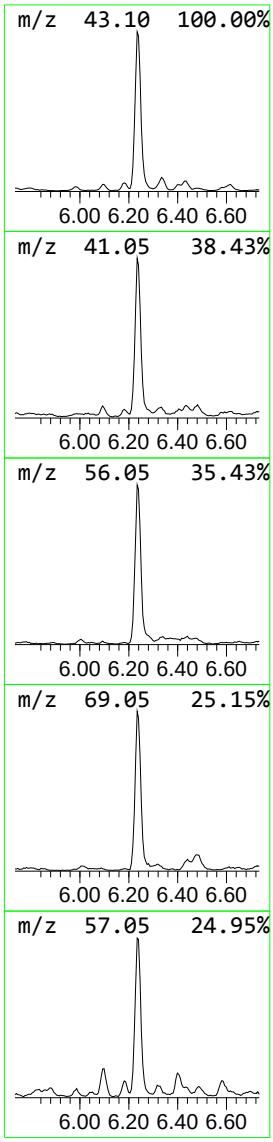
****

Peak Number 1 2-Methyl-1-hexanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.234	32.53 ng/ul	1713100	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-hexanol	116	C7H16O	1000427-67-0	83
2	Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	64
3	Hexane, 1-propoxy-	144	C9H20O	053685-78-2	59
4	1-Pentanol, 2-ethyl-	116	C7H16O	027522-11-8	53
5	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

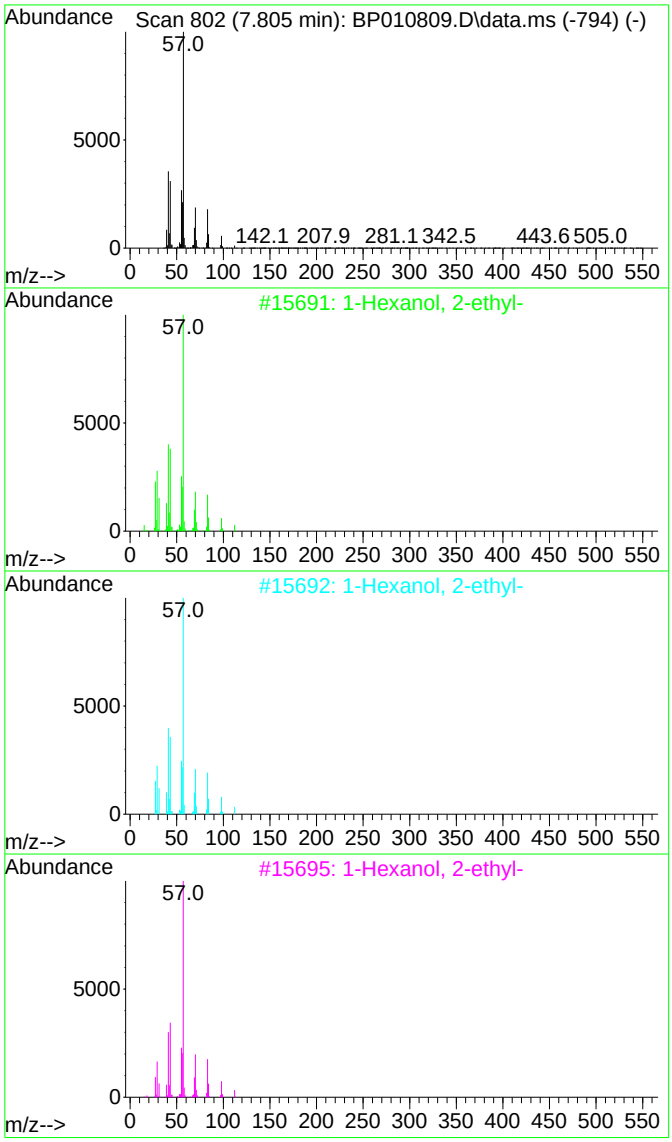
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

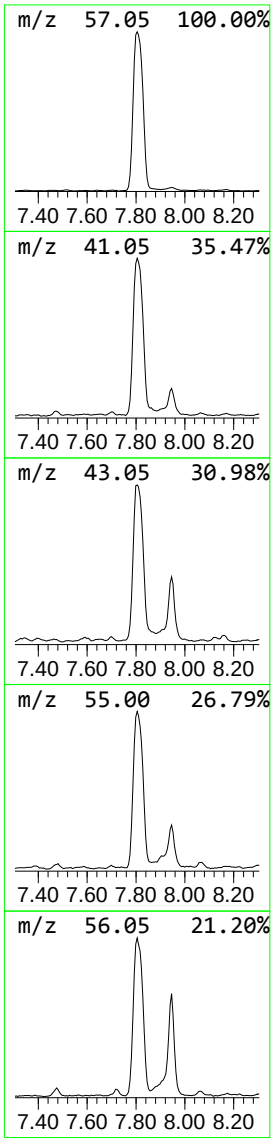
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.805	99.19 ng/ul	5222940	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
5	Oxalic acid, isobutyl octyl ester	258	C14H26O4	1000309-37-3	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

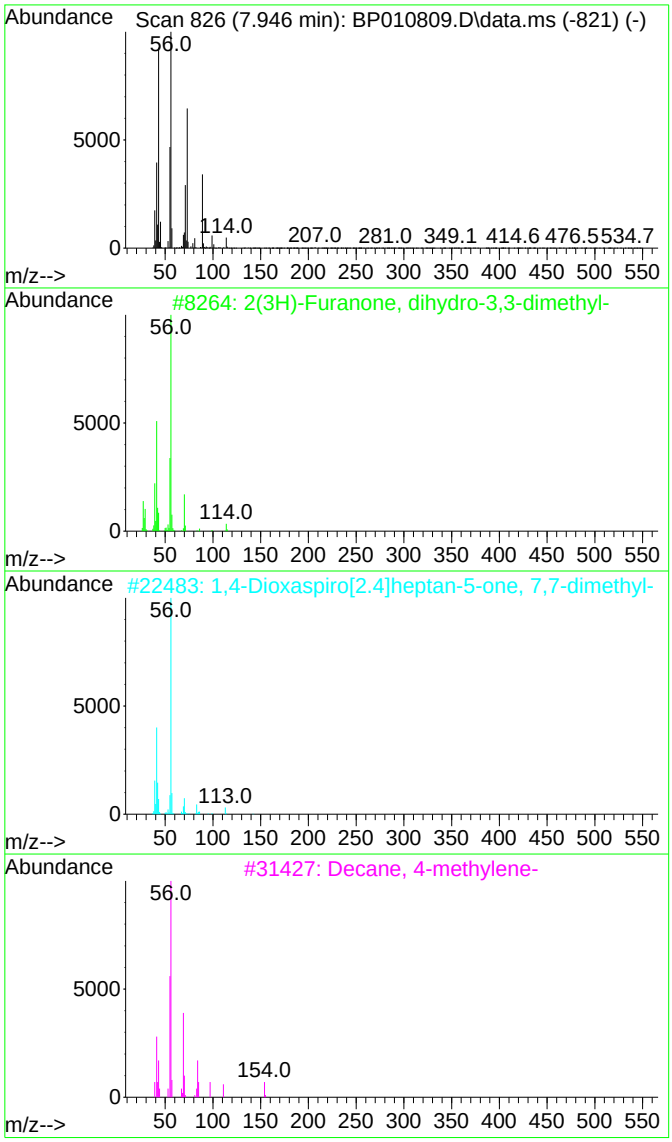
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

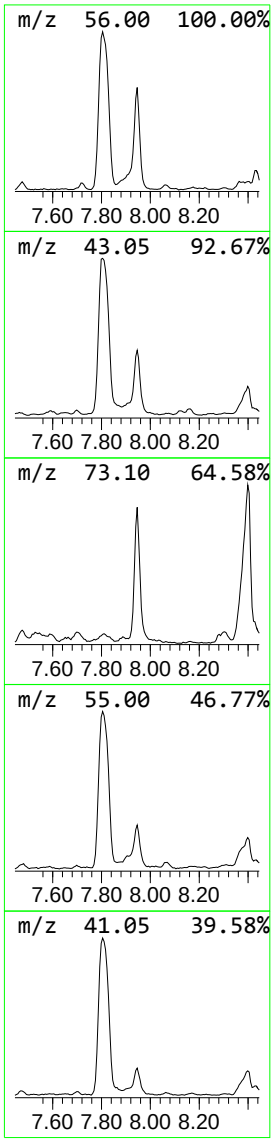
****

Peak Number 3 unknown-01 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.946	12.57 ng/ul	661702	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Furanone, dihydro-3,3-dime...	114	C6H10O2	003709-08-8	47
2	1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	37
3	Decane, 4-methylene-	154	C11H22	024949-41-5	37
4	2,5-Dimethyl-5-hexen-3-ol	128	C8H16O	067760-91-2	33
5	[1,4]Dioxino[2,3-b]-1,4-dioxin, ...	202	C10H18O4	038737-48-3	33





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

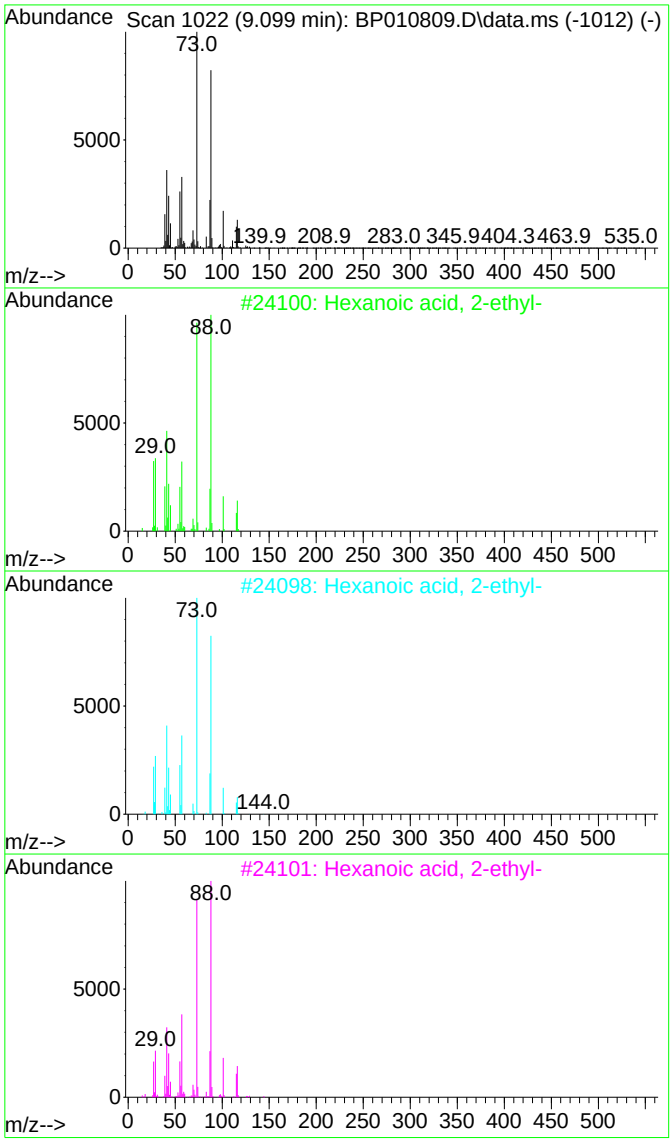
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

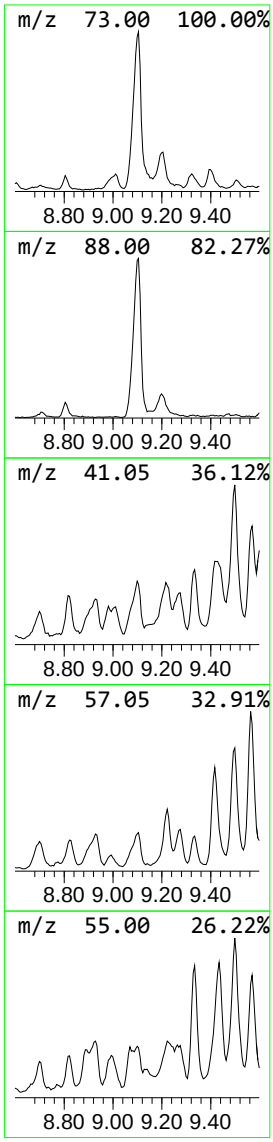
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.099	18.90 ng/ul	995214	1,4-Dichlorobenzene-d4	7.722

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

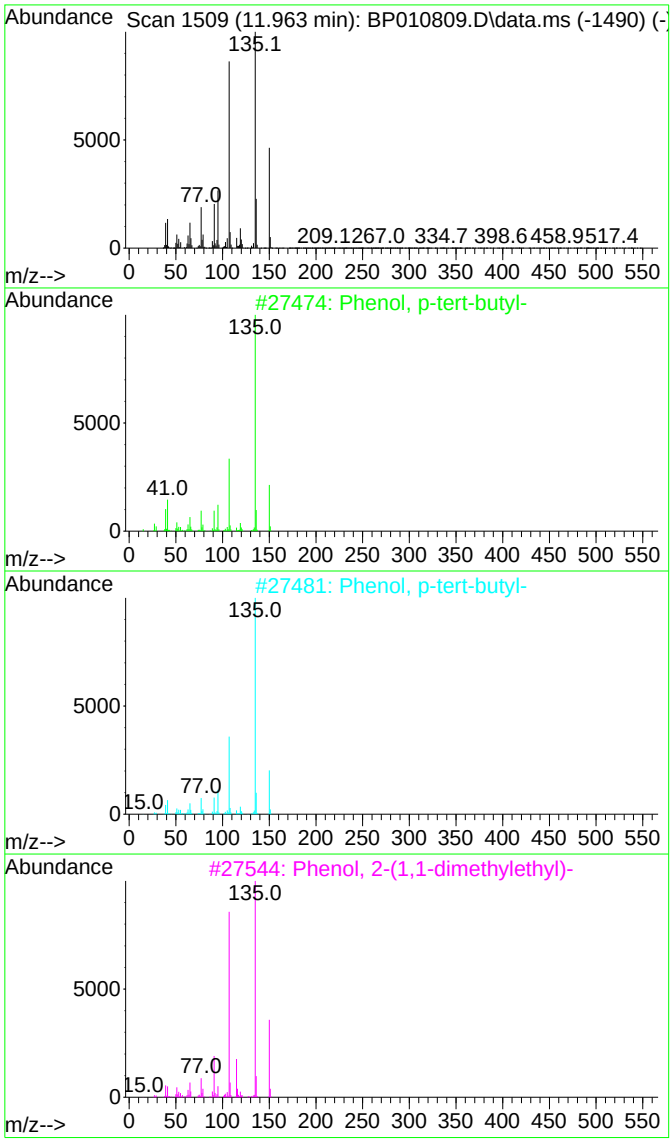
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

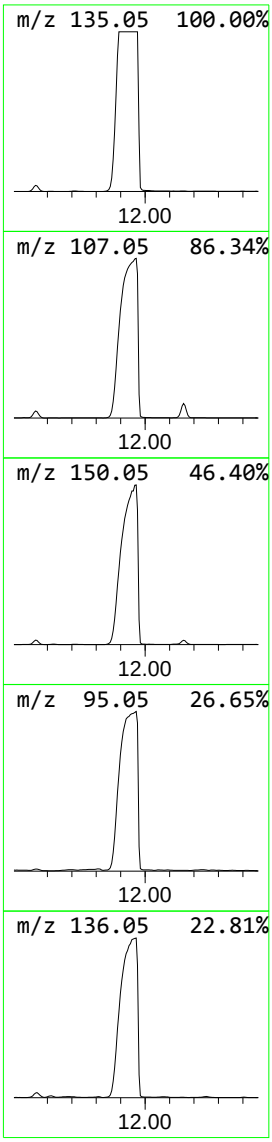
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	21.18 ng/ul	230514000	Naphthalene-d8	10.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
3	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	80
4	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	72
5	1,3-Cyclopentadiene, 1,3-bis(1-m...	150	C11H18	123278-27-3	64





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Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

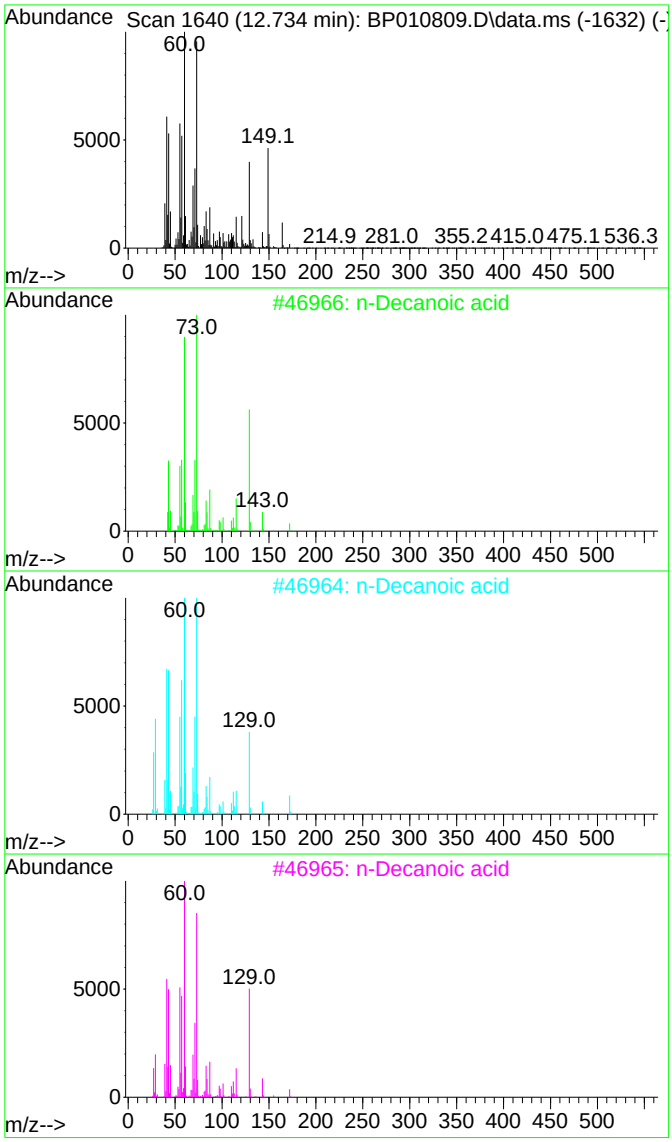
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

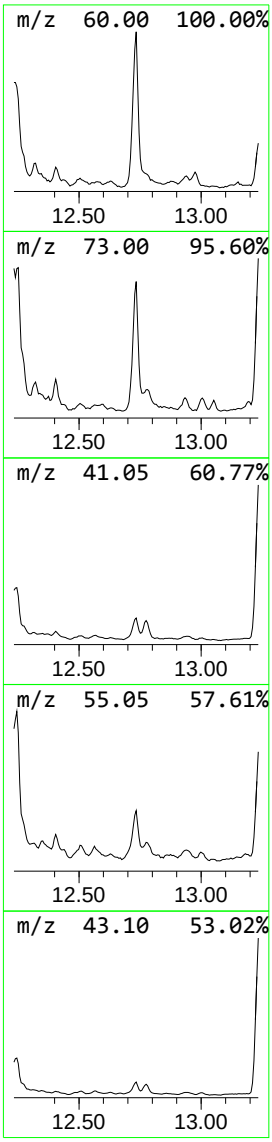
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 n-Decanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.734	25.38 ng/ul	2326440	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	95
2	n-Decanoic acid	172	C10H20O2	000334-48-5	89
3	n-Decanoic acid	172	C10H20O2	000334-48-5	68
4	n-Decanoic acid	172	C10H20O2	000334-48-5	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	59





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

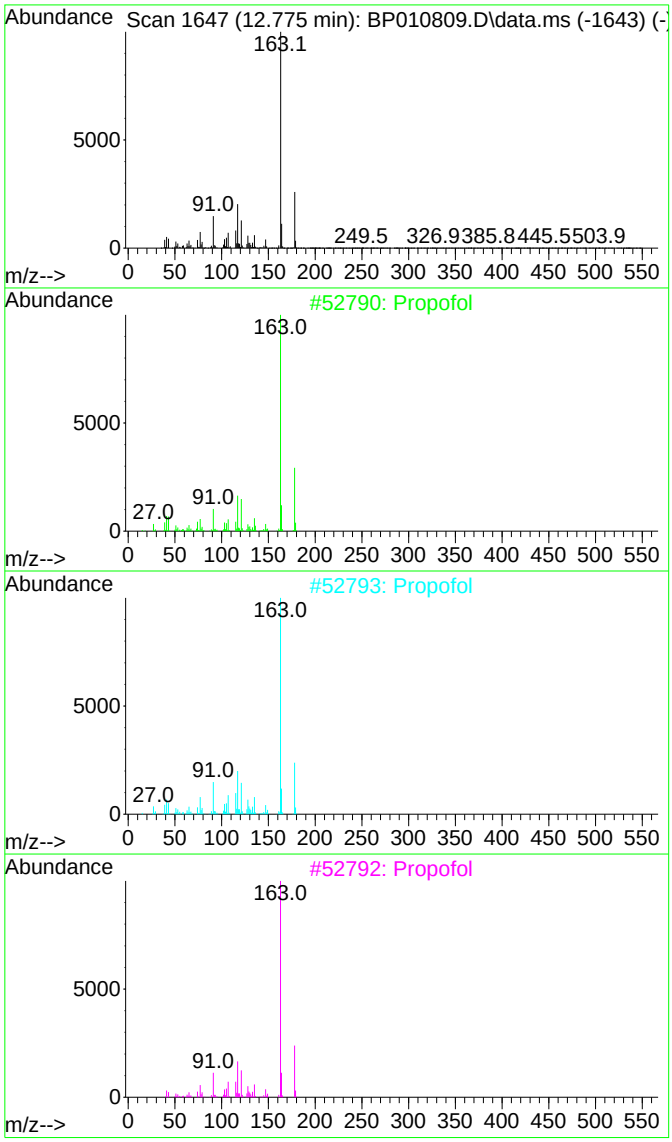
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

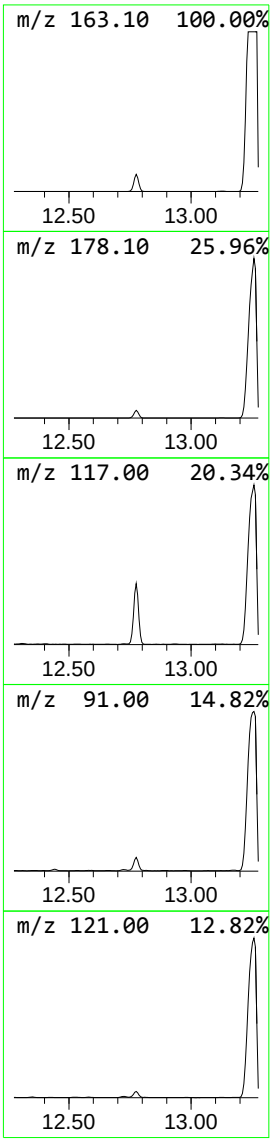
****

Peak Number 9 Propofol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.775	52.34 ng/ul	4798080	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	97
2	Propofol	178	C12H18O	002078-54-8	97
3	Propofol	178	C12H18O	002078-54-8	96
4	Propofol	178	C12H18O	002078-54-8	94
5	Propofol	178	C12H18O	002078-54-8	94





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

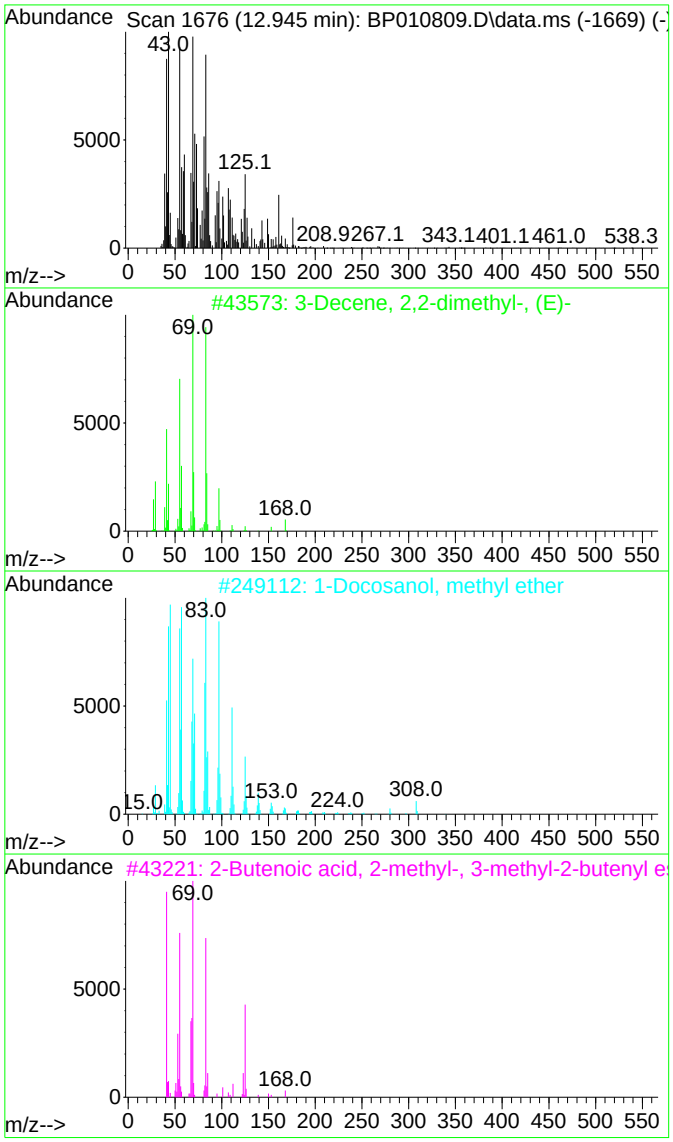
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

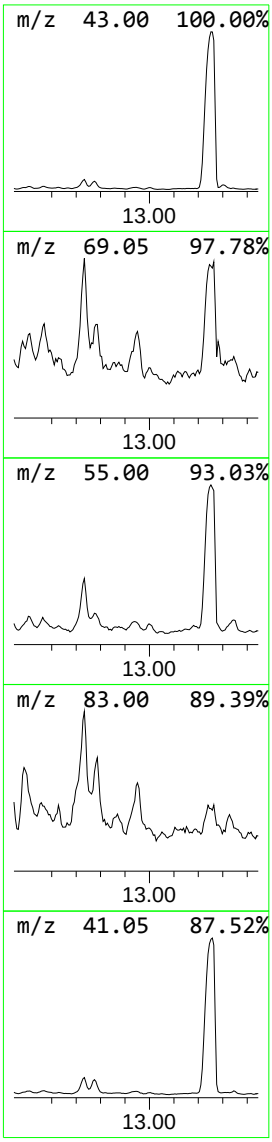
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.945	5.75 ng/ul	527525	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Decene, 2,2-dimethyl-, (E)-	168	C12H24	055499-02-0	43
2	1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	38
3	2-Butenoic acid, 2-methyl-, 3-me...	168	C10H16O2	083783-86-2	30
4	2-Undecene, 8-methyl-, (Z)-	168	C12H24	074630-44-7	25
5	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	15





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

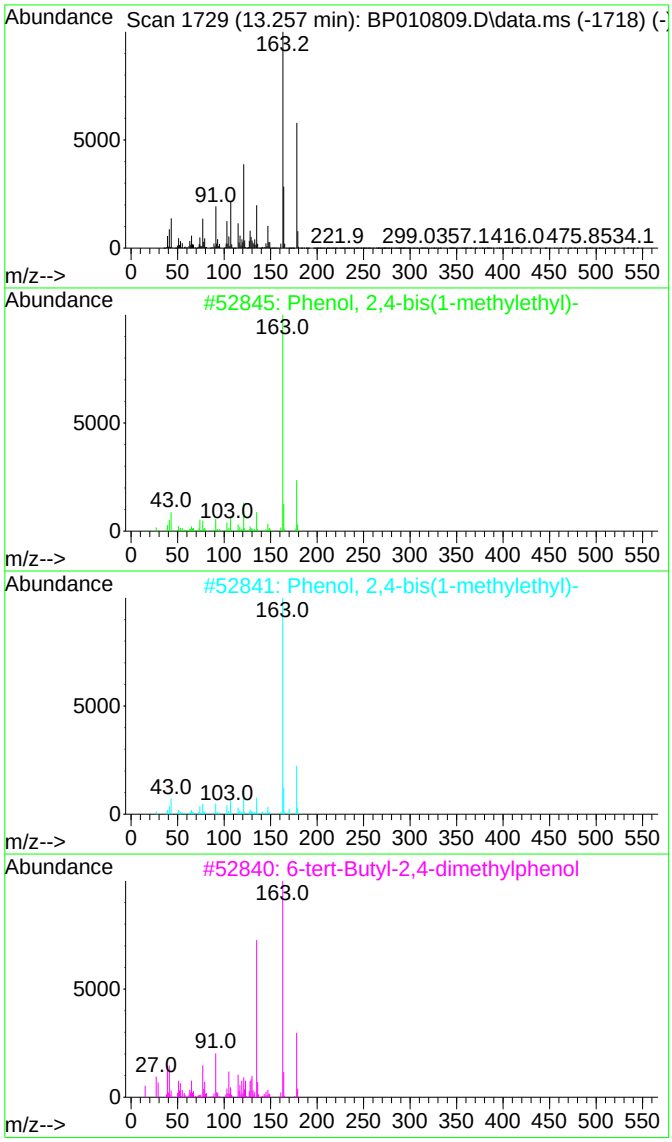
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

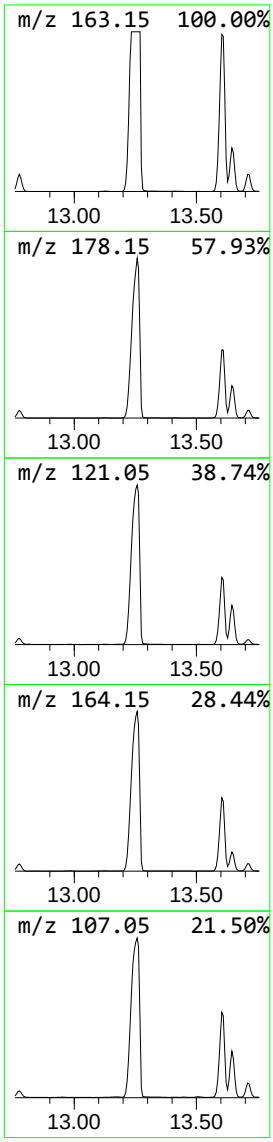
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Phenol, 2,4-bis(1-methyleth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.257	1285.91 ng/ul	117881000	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90
2	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	78
3	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	76
4	Propofol	178	C12H18O	002078-54-8	72
5	Phenol, 2-(1,1-dimethylethyl)-4-...	178	C12H18O	000096-70-8	64





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

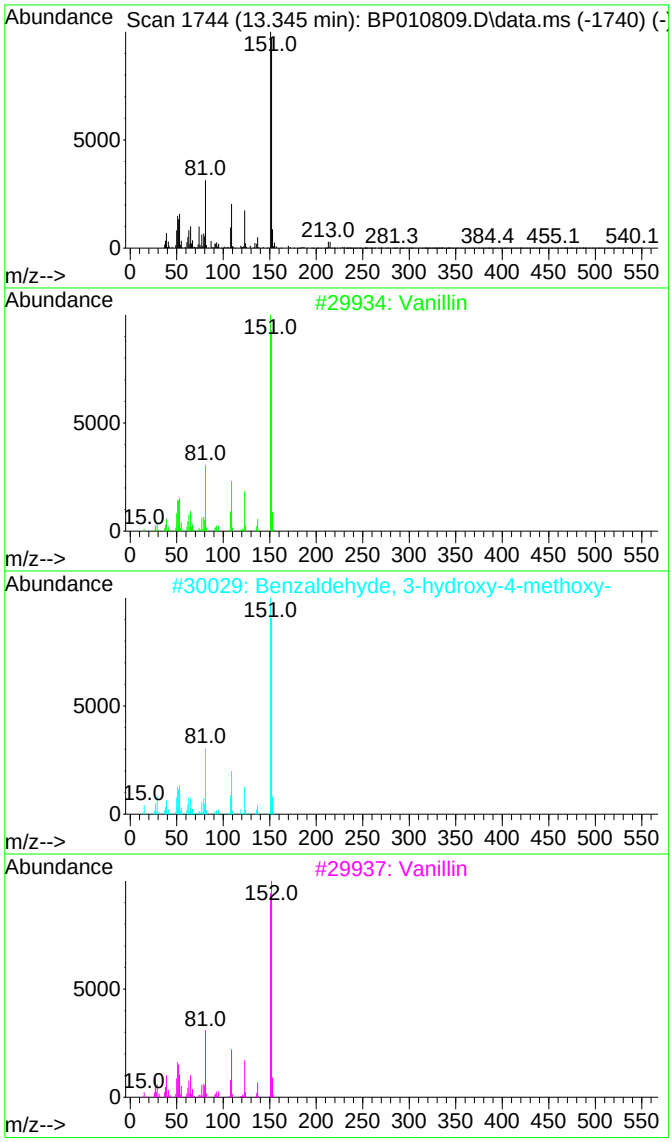
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

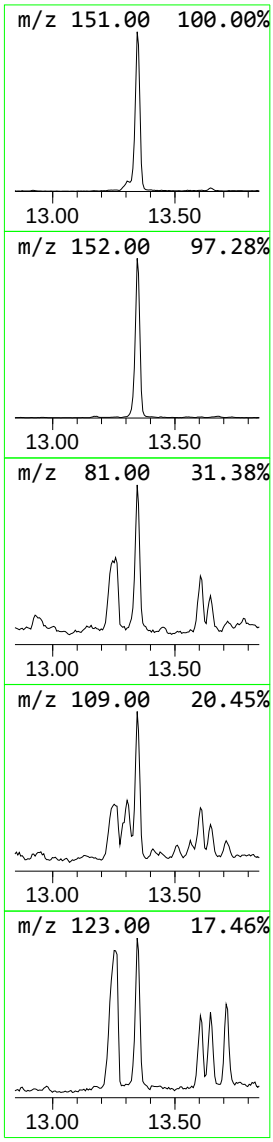
****

Peak Number 12 Vanillin Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.345	21.18 ng/ul	1941260	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	97
2	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	97
3	Vanillin	152	C8H8O3	000121-33-5	97
4	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96
5	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

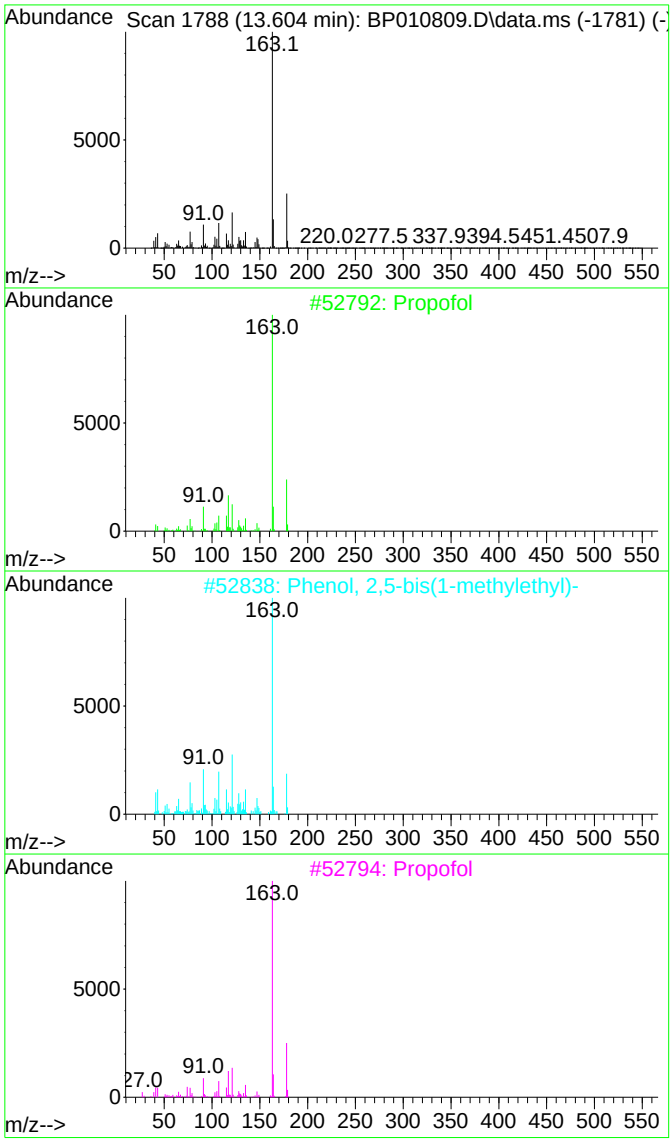
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

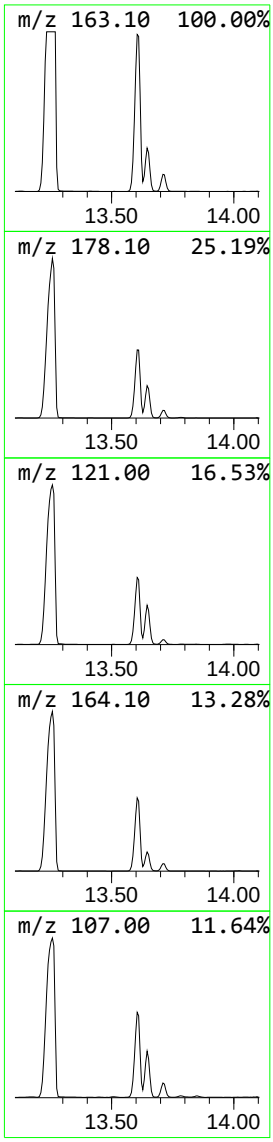
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Phenol, 2,5-bis(1-methyleth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.604	442.26 ng/ul	40542700	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	97
2	Phenol, 2,5-bis(1-methylethyl)-	178	C12H18O	035946-91-9	95
3	Propofol	178	C12H18O	002078-54-8	94
4	Propofol	178	C12H18O	002078-54-8	94
5	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90





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Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

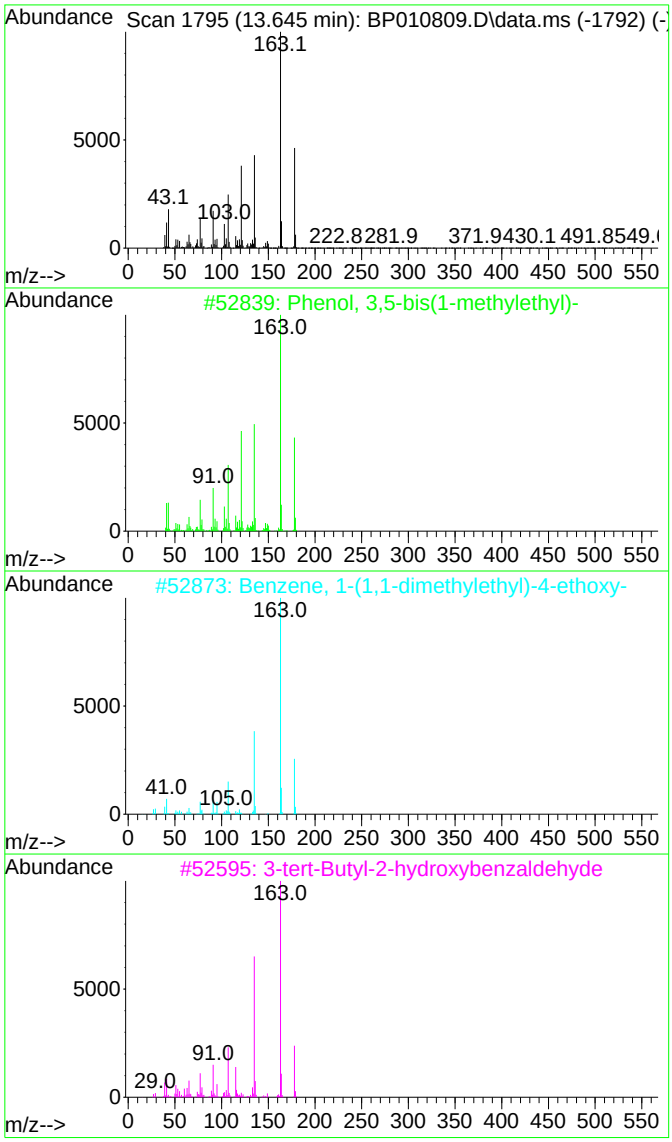
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

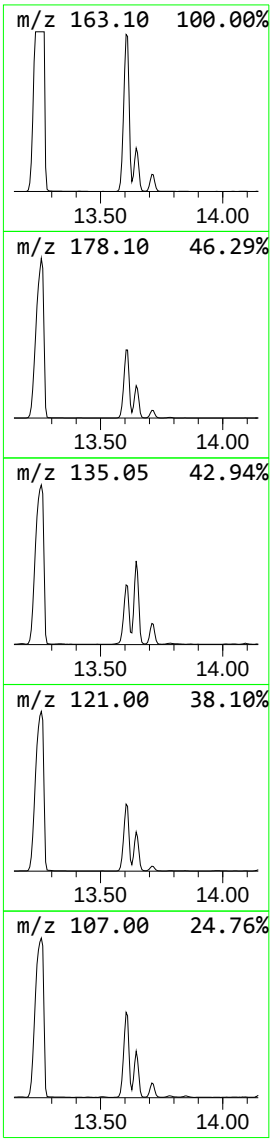
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Phenol, 3,5-bis(1-methyleth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.645	153.06 ng/ul	14031400	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-bis(1-methylethyl)-	178	C12H18O	026886-05-5	97
2	Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	90
3	3-tert-Butyl-2-hydroxybenzaldehyde	178	C11H14O2	024623-65-2	86
4	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	81
5	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	81





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

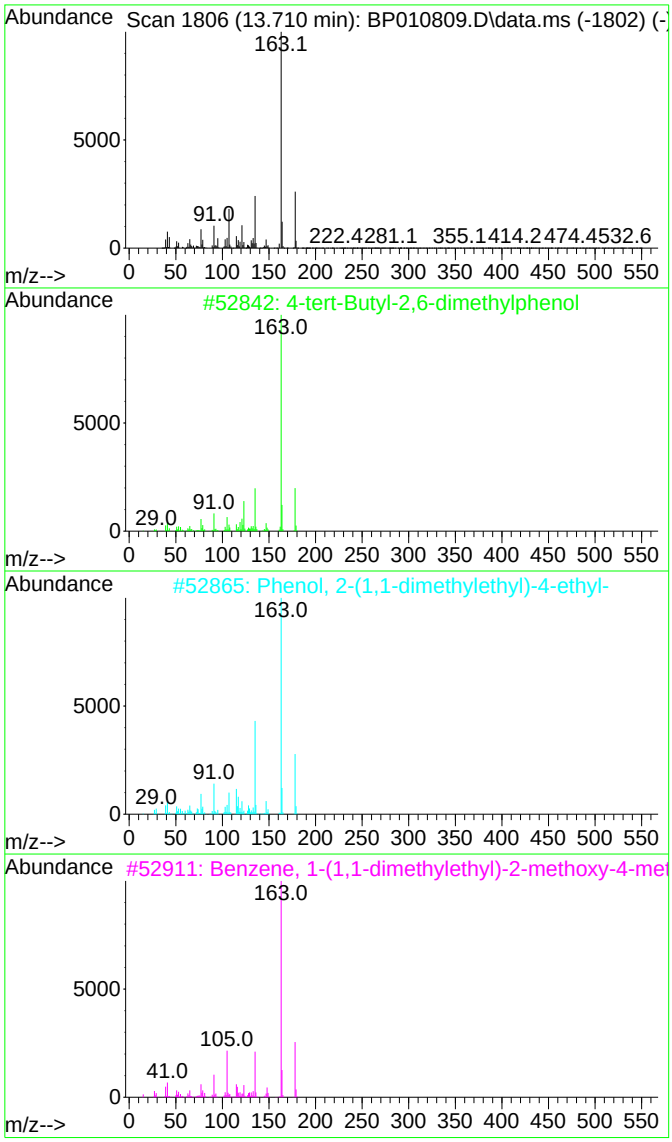
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

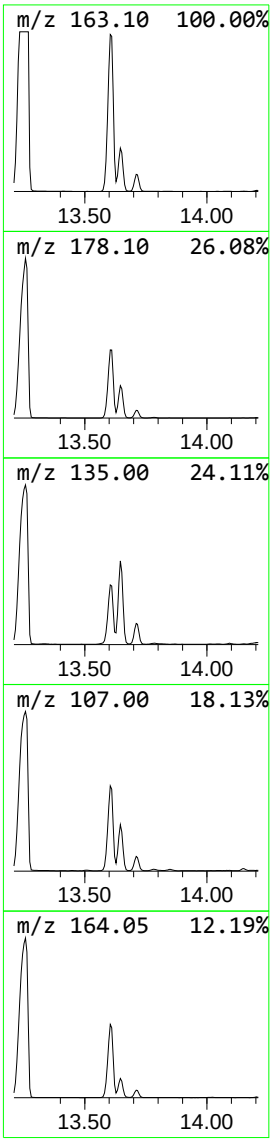
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 4-tert-Butyl-2,6-dimethylph... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.710	47.60 ng/ul	4363520	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-tert-Butyl-2,6-dimethylphenol	178	C12H18O	000879-97-0	91
2	Phenol, 2-(1,1-dimethylethyl)-4-...	178	C12H18O	000096-70-8	87
3	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	83
4	Propofol	178	C12H18O	002078-54-8	83
5	Propofol	178	C12H18O	002078-54-8	81





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

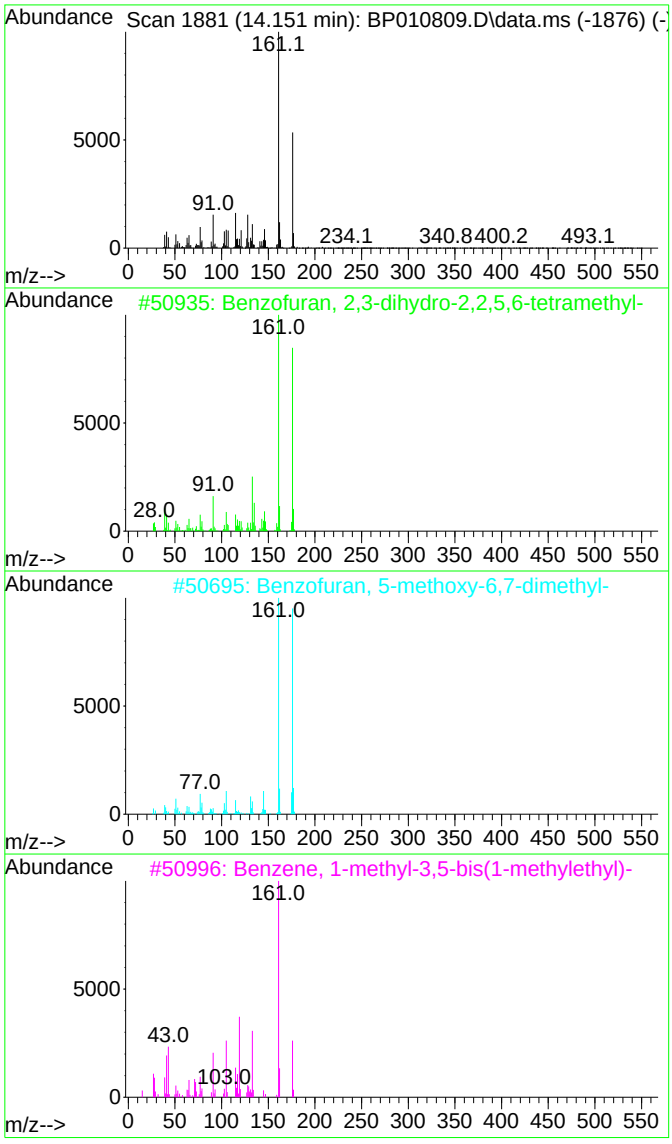
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

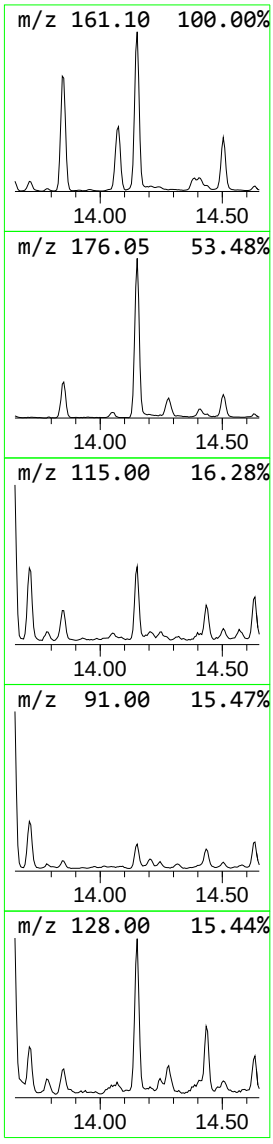
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzofuran, 2,3-dihydro-2,2... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.151	19.08 ng/ul	1748990	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2,3-dihydro-2,2,5,6-...	176	C12H16O	063577-97-9	90
2	Benzofuran, 5-methoxy-6,7-dimethyl-	176	C11H12O2	035355-35-2	80
3	Benzene, 1-methyl-3,5-bis(1-meth...	176	C13H20	003055-14-9	80
4	Benzene, 1,2,3,4-tetramethyl-5-(...	176	C13H20	061142-67-4	72
5	(1R)-1-(4-tert-butylphenyl)ethan...	177	C12H19N	089538-65-8	72





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Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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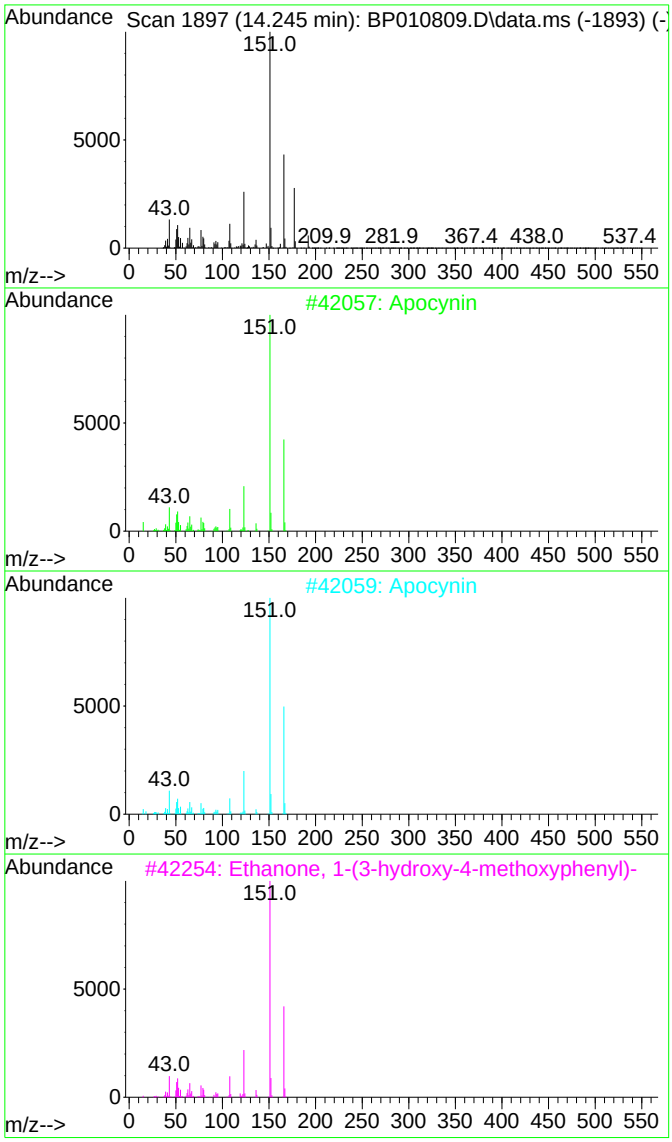
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

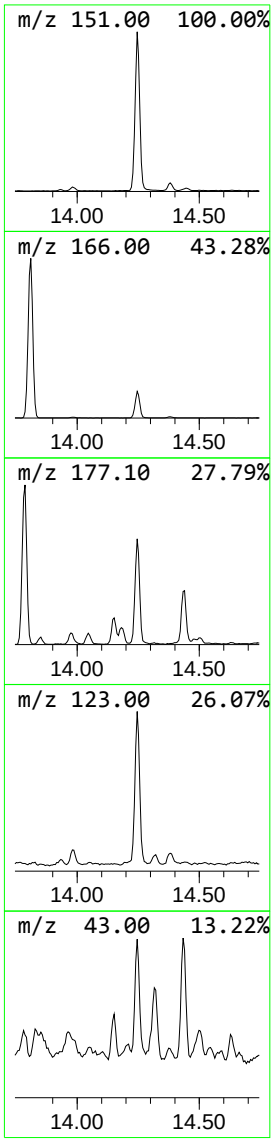
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Apocynin Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.245	17.68 ng/ul	1620590	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Apocynin	166	C9H10O3	000498-02-2	96
2	Apocynin	166	C9H10O3	000498-02-2	95
3	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	95
4	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	93
5	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	91





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Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

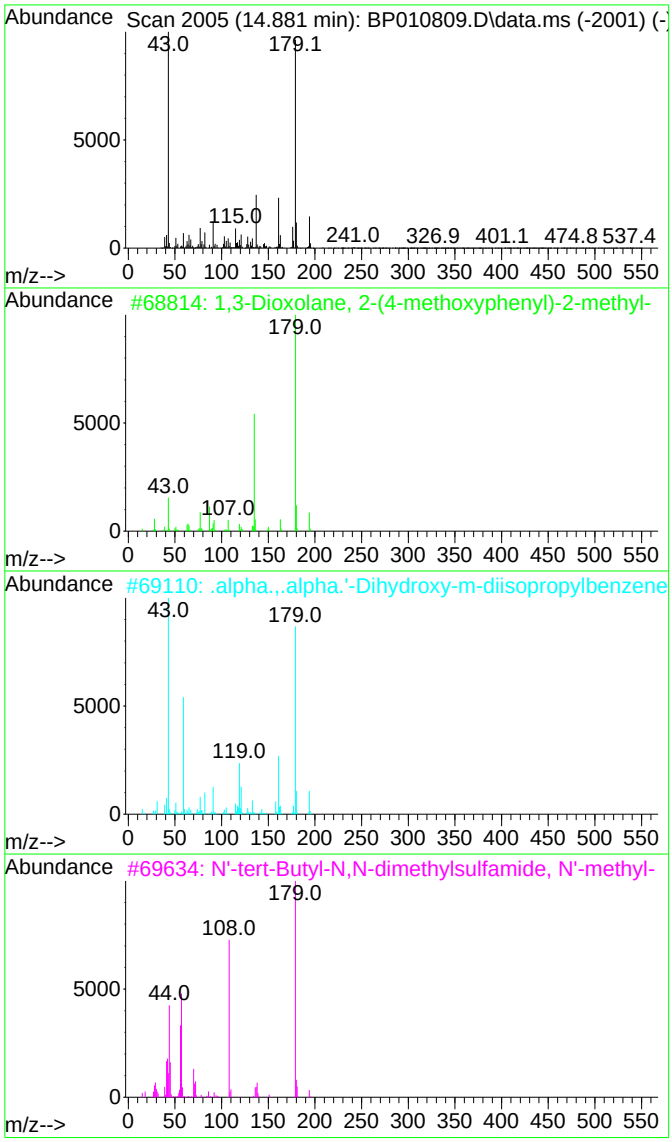
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

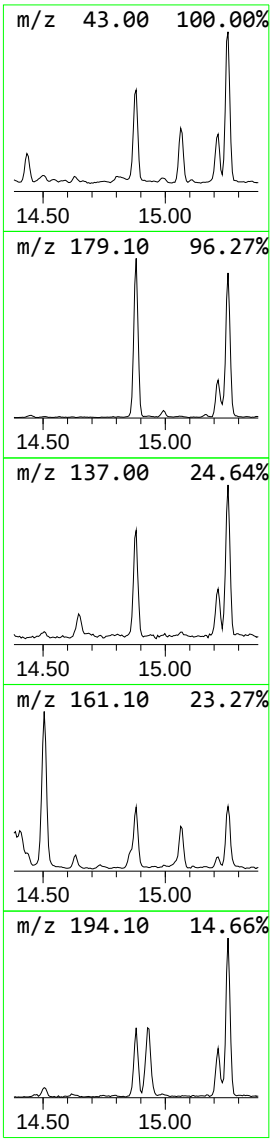
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 1,3-Dioxolane, 2-(4-methoxy... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.881	9.54 ng/ul	874570	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 2-(4-methoxypheny...	194	C11H14O3	036881-00-2	50
2	.alpha.,.alpha.'-Dihydroxy-m-dii...	194	C12H18O2	001999-85-5	50
3	N'-tert-Butyl-N,N-dimethylsulfam...	194	C7H18N2O2S	1858529-07-3	50
4	5-tert-Butyl-4-hydroxymethyl-2-m...	212	C11H16O4	1000275-36-7	47
5	1,4-Bis[.alpha.-methoxyethyl]ben...	194	C12H18O2	129770-42-9	47





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Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

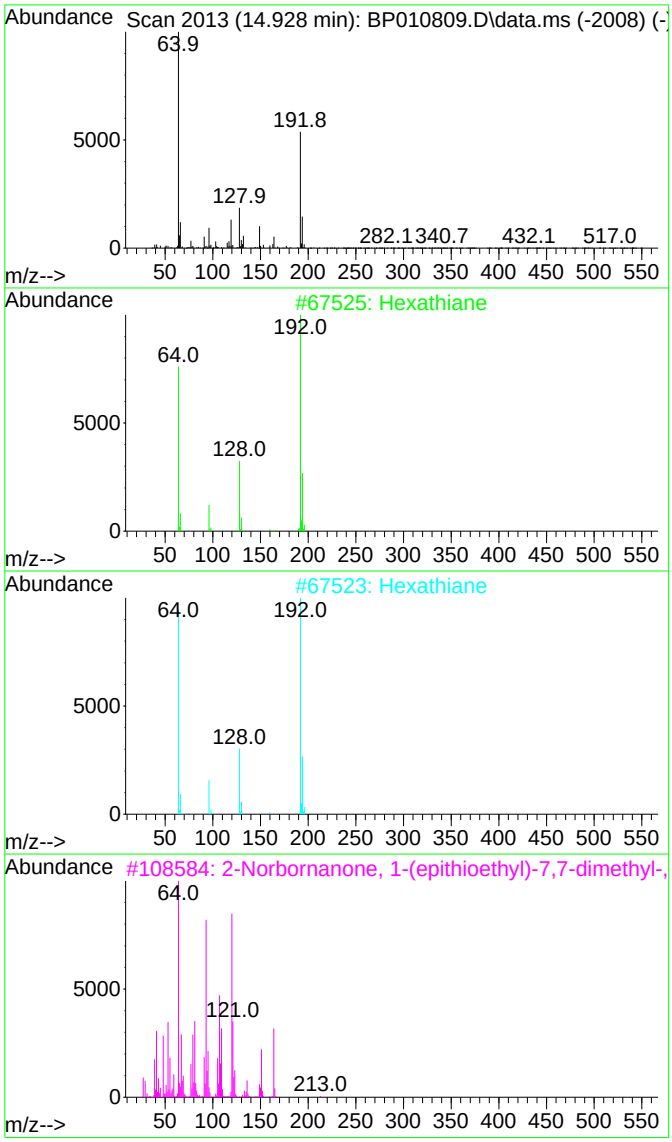
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

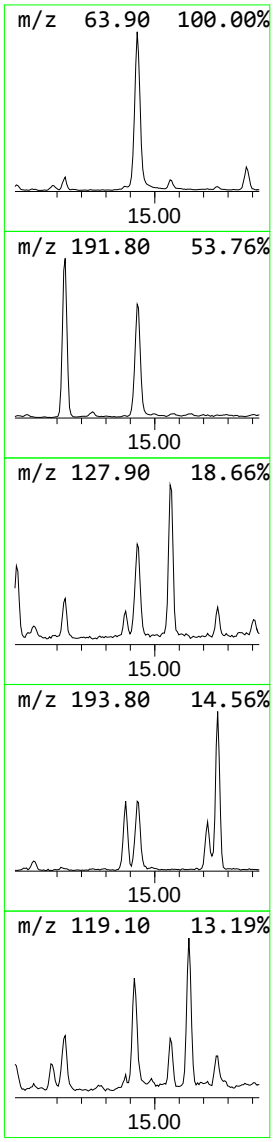
****

Peak Number 20 Hexathiane Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.928	9.38 ng/ul	859679	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexathiane	192	S6	013798-23-7	91
2	Hexathiane	192	S6	013798-23-7	86
3	2-Norbornanone, 1-(epithioethyl)...	228	C11H16O3S	001782-93-0	59
4	Di(tert-butyl) trisulfide, perfl...	534	C8F18S3	016005-64-4	50
5	Hexathiane	192	S6	013798-23-7	11





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

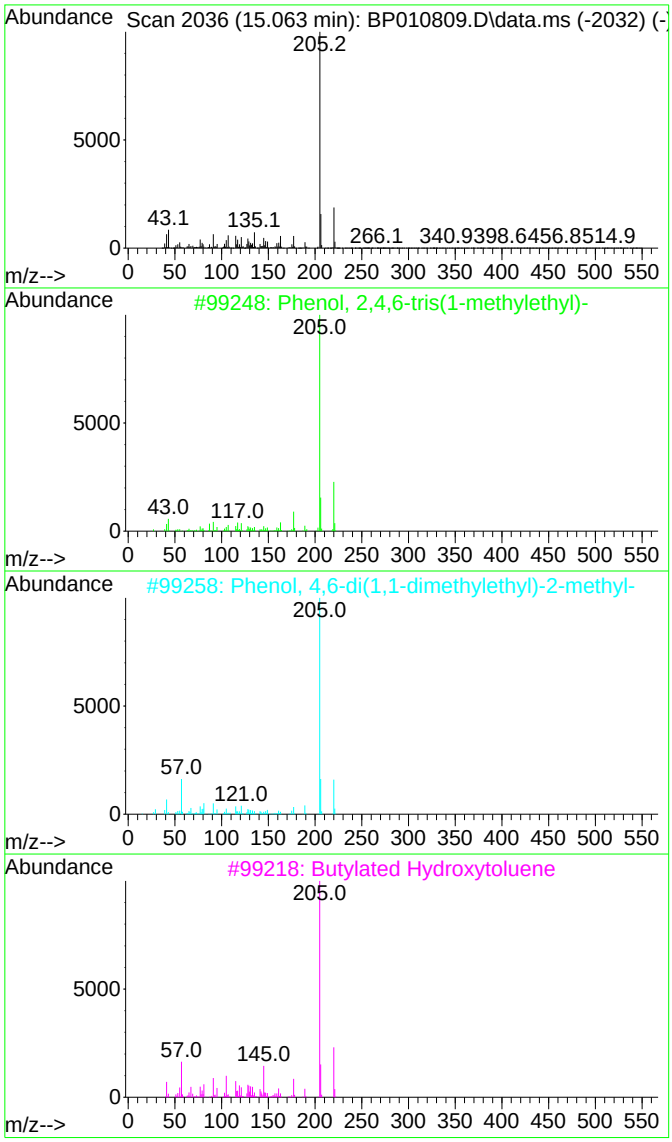
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

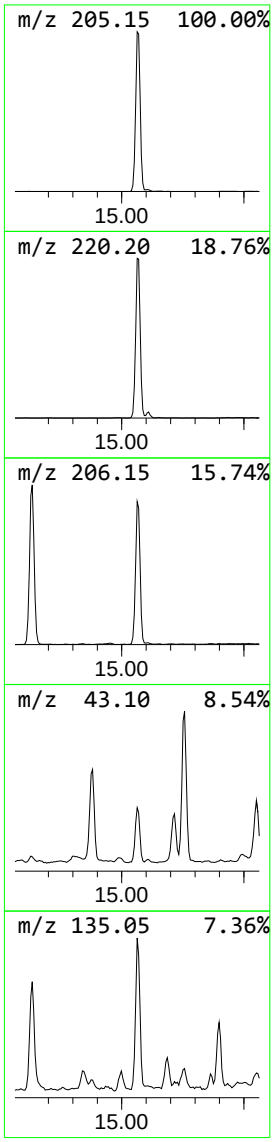
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Phenol, 2,4,6-tris(1-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.063	44.72 ng/ul	4099650	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	87
2	Phenol, 4,6-di(1,1-dimethylethyl...	220	C15H24O	000616-55-7	83
3	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	81
4	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	74
5	Phenol, 4,6-di(1,1-dimethylethyl...	220	C15H24O	000616-55-7	74





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

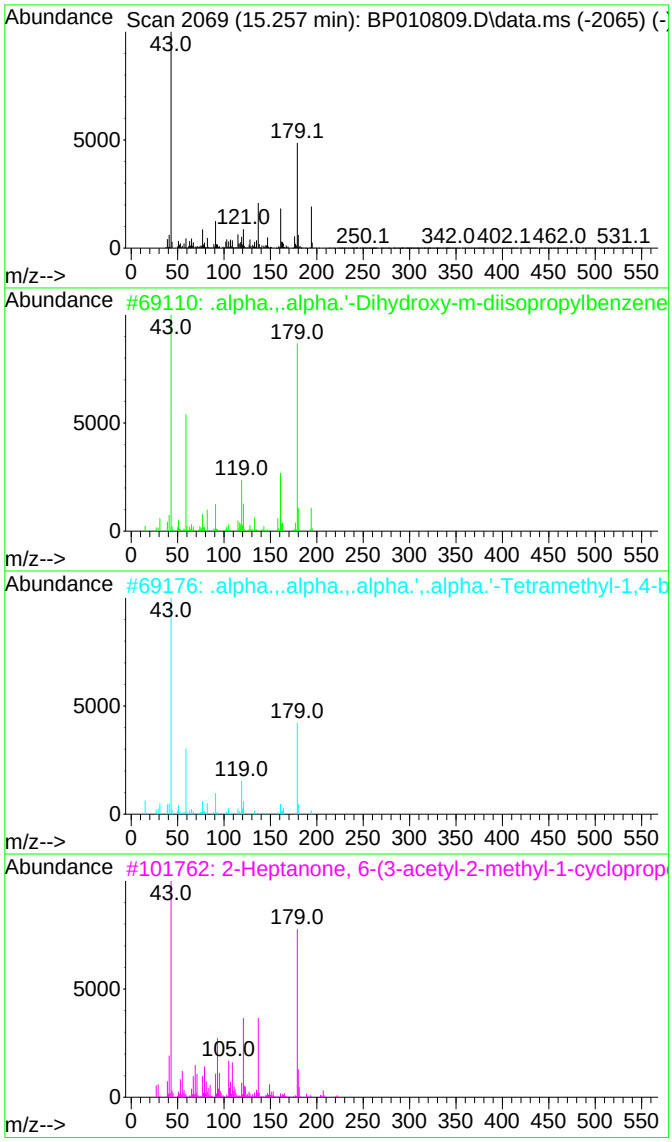
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

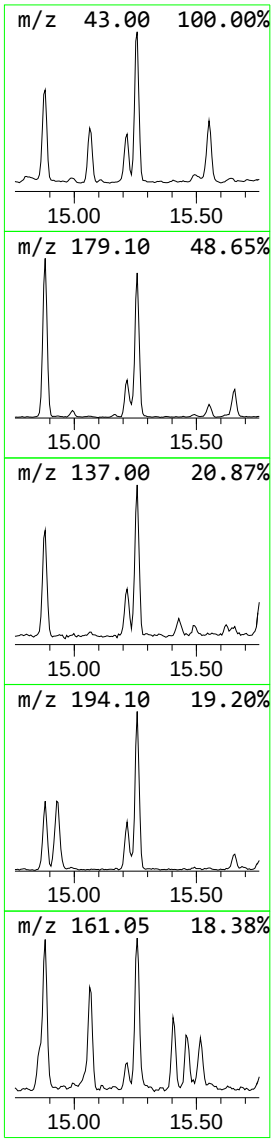
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 unknown-02 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.257	12.48 ng/ul	1143630	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.,.alpha.'-Dihydroxy-m-dii...	194	C12H18O2	001999-85-5	37
2	.alpha.,.alpha.,.alpha.',.alpha....	194	C12H18O2	002948-46-1	25
3	2-Heptanone, 6-(3-acetyl-2-methy...	222	C14H22O2	065868-86-2	9
4	2,3-Dimethoxyphenethylamine	181	C10H15NO2	003213-29-4	7
5	Ethanone, 1-(1-hydroxy-2,6,6-tri...	180	C11H16O2	058254-18-5	7





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

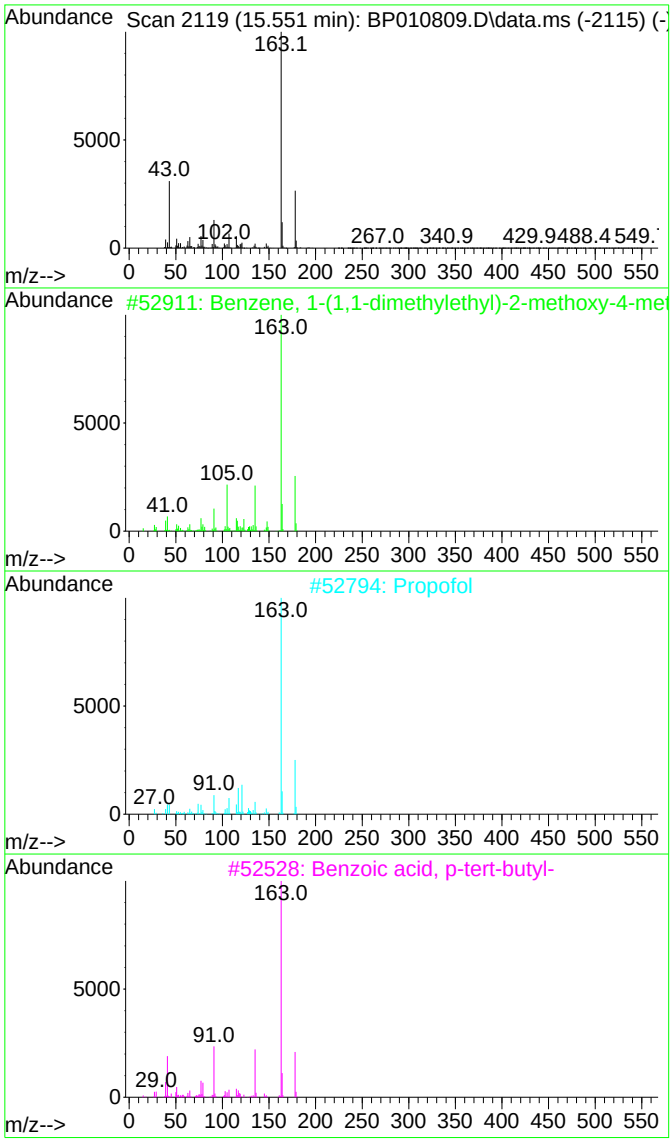
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

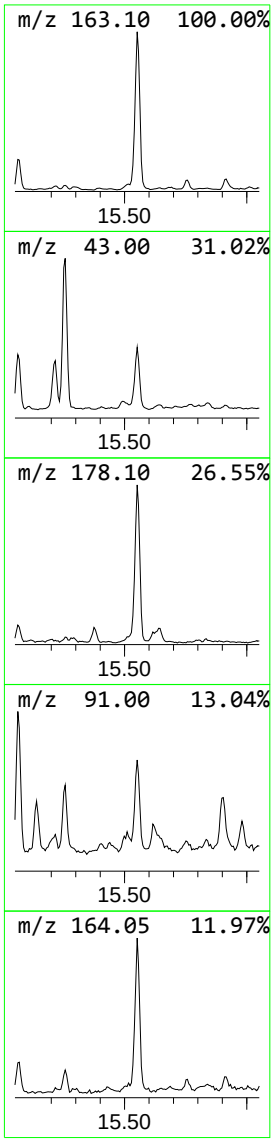
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1-(1,1-dimethyleth... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.551	14.11 ng/ul	1293370	Acenaphthene-d10	14.381

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	86
2	Propofol	178	C12H18O	002078-54-8	86
3	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	80
4	Propofol	178	C12H18O	002078-54-8	78
5	Propofol	178	C12H18O	002078-54-8	78





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Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

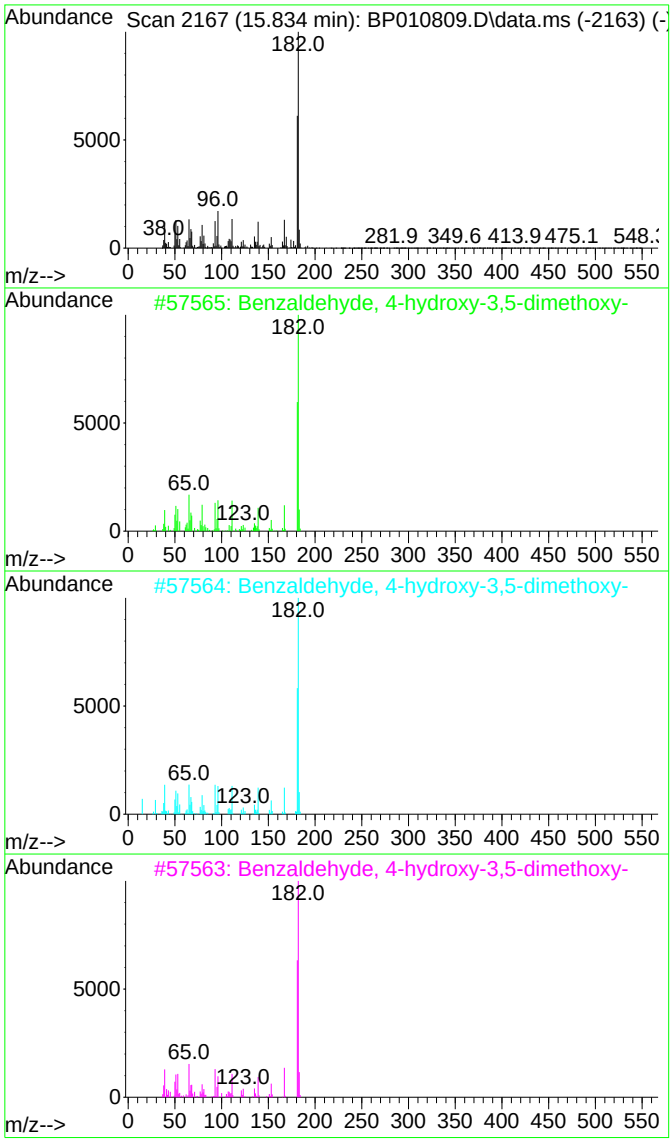
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

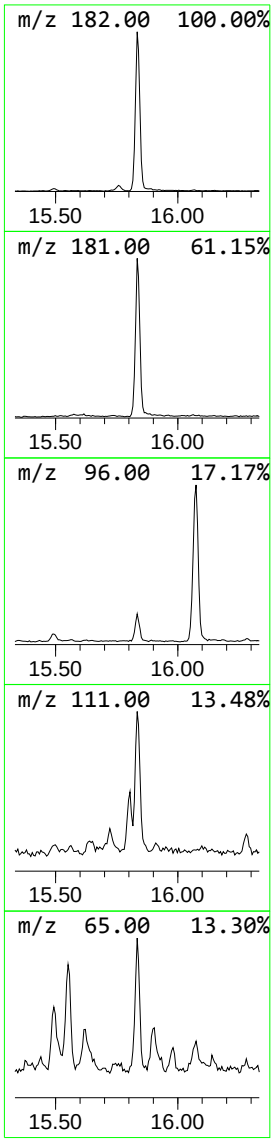
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzaldehyde, 4-hydroxy-3,5... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.834	11.13 ng/ul	1096880	Phenanthrene-d10	17.139

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	97
2	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	97
3	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	97
4	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96
5	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96





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Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

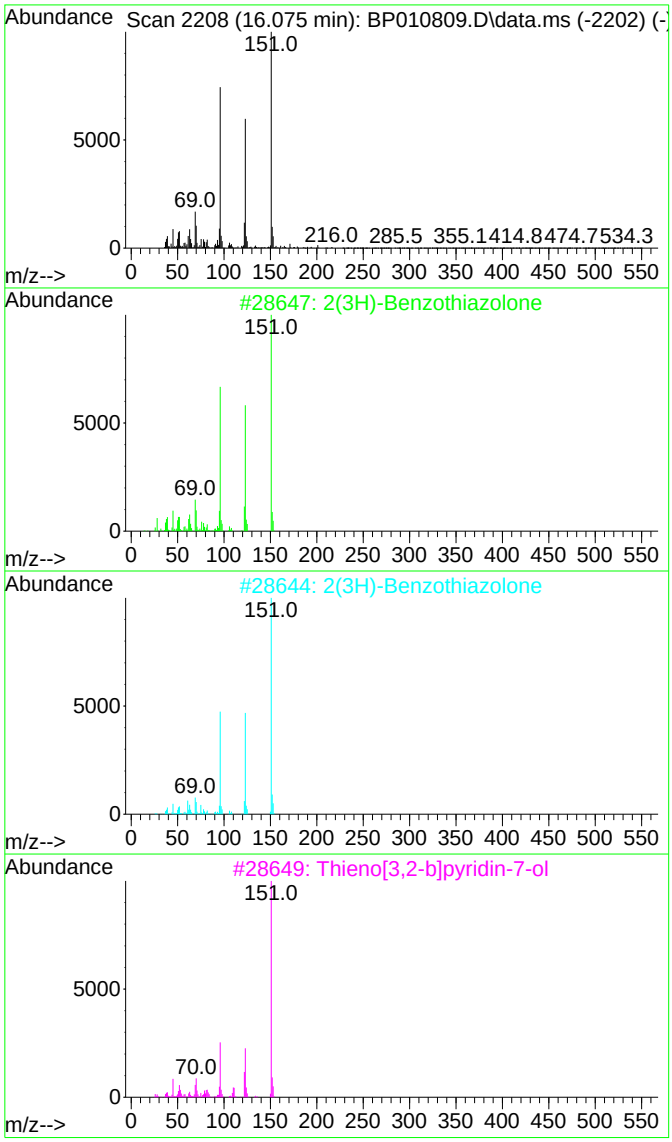
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

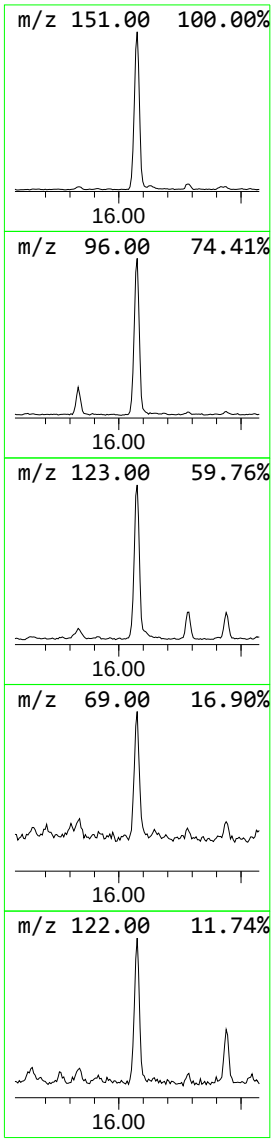
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 2(3H)-Benzothiazolone Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	16.62 ng/ul	1637260	Phenanthrene-d10	17.139

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	97
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	93
3	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	58
4	t-Butylhydroquinone	166	C10H14O2	001948-33-0	43
5	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	43





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

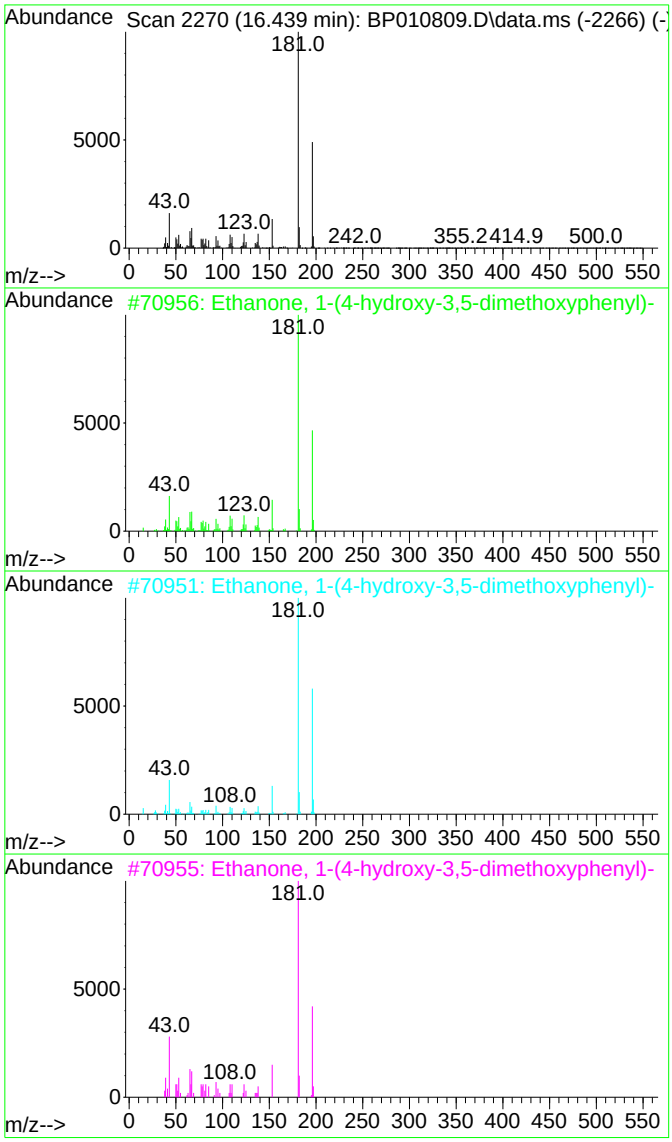
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

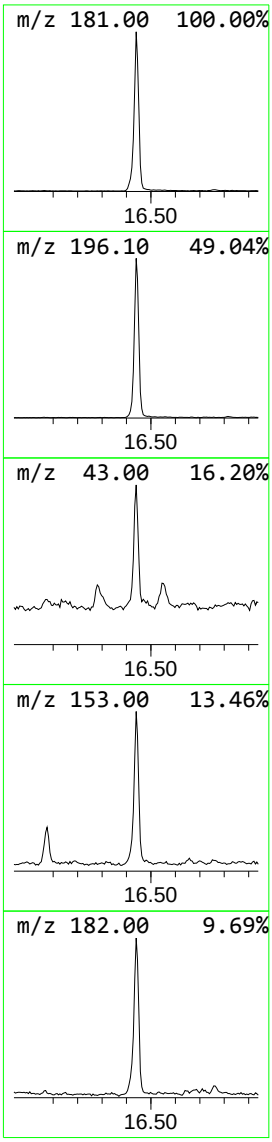
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Ethanone, 1-(4-hydroxy-3,5-... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.439	13.40 ng/ul	1319920	Phenanthrene-d10	17.139

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	98
2	Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	97
3	Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	96
4	Ethanone, 1-(4-hydroxy-3,5-dimet...	196	C10H12O4	002478-38-8	96
5	4,5-Dimethoxy-2-hydroxyacetophenone	196	C10H12O4	020628-06-2	90





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Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

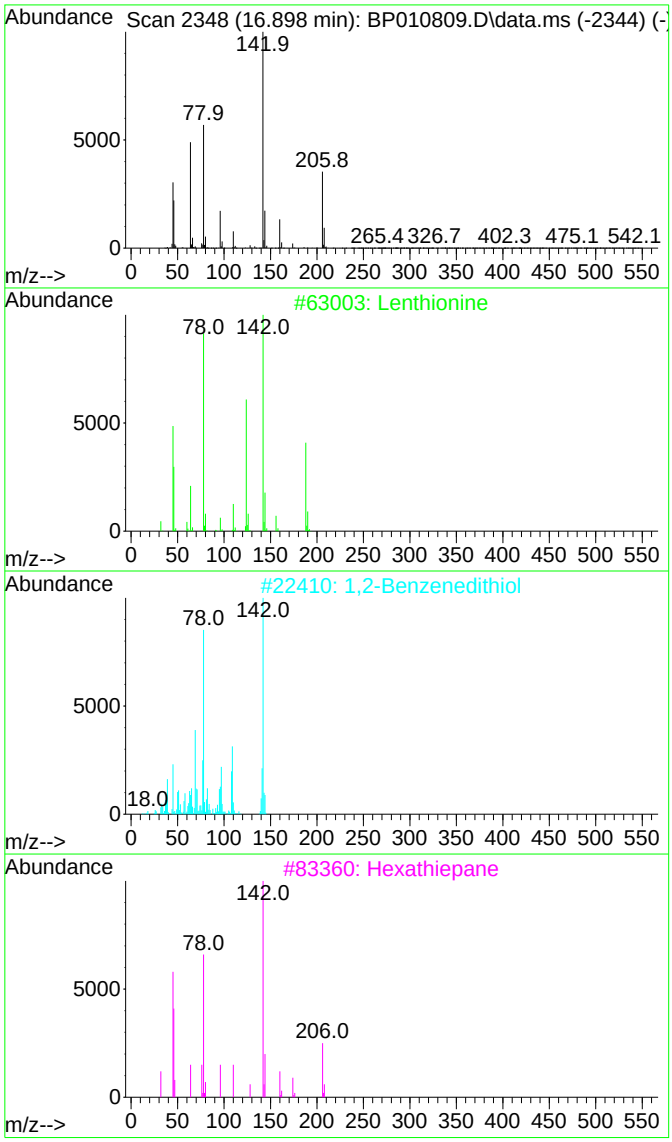
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

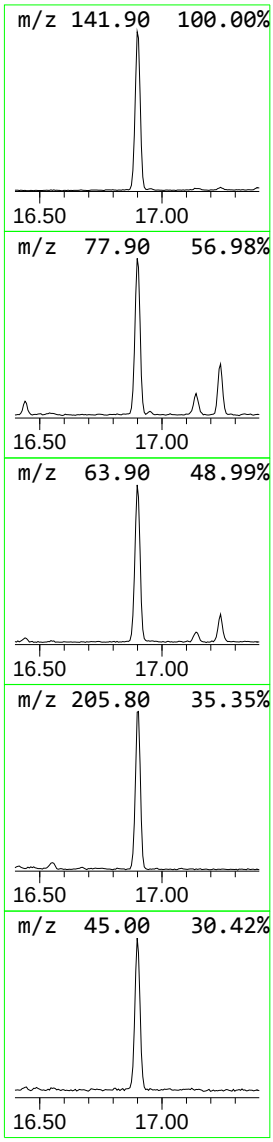
****

Peak Number 27 unknown-03 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.898	13.46 ng/ul	1326500	Phenanthrene-d10	17.139

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lenthionine	188	C2H4S5	000292-46-6	22
2	1,2-Benzenedithiol	142	C6H6S2	017534-15-5	12
3	Hexathiepane	206	CH2S6	017233-71-5	12
4	4(1H)-Pyrimidinone, 2-(methylthio)-	142	C5H6N2OS	005751-20-2	9
5	3,4-Ethylenedioxythiophene	142	C6H6O2S	126213-50-1	9





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Sample : N3374-01
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ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

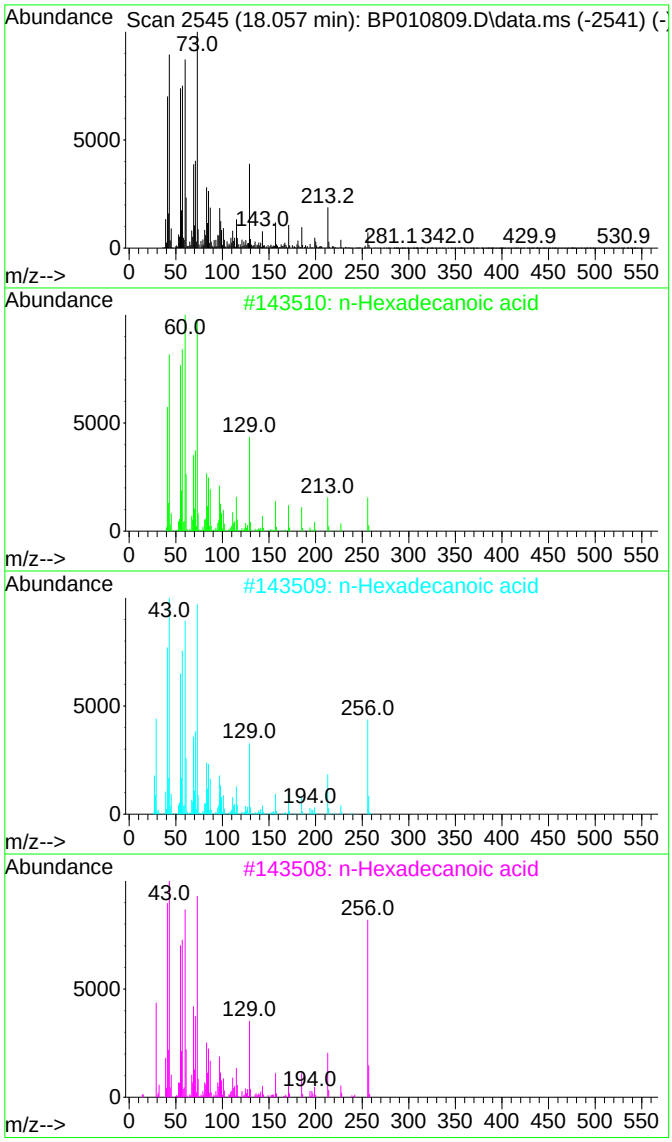
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

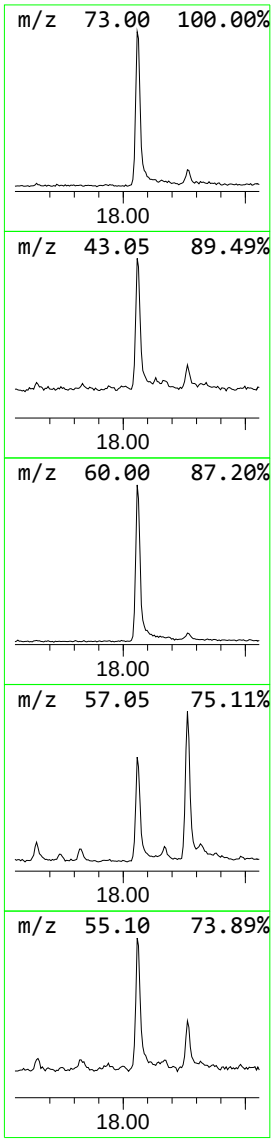
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 n-Hexadecanoic acid Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.057	15.72 ng/ul	1548920	Phenanthrene-d10	17.139

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





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ALS Vial : 28 Sample Multiplier: 1

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ClientSampleId :
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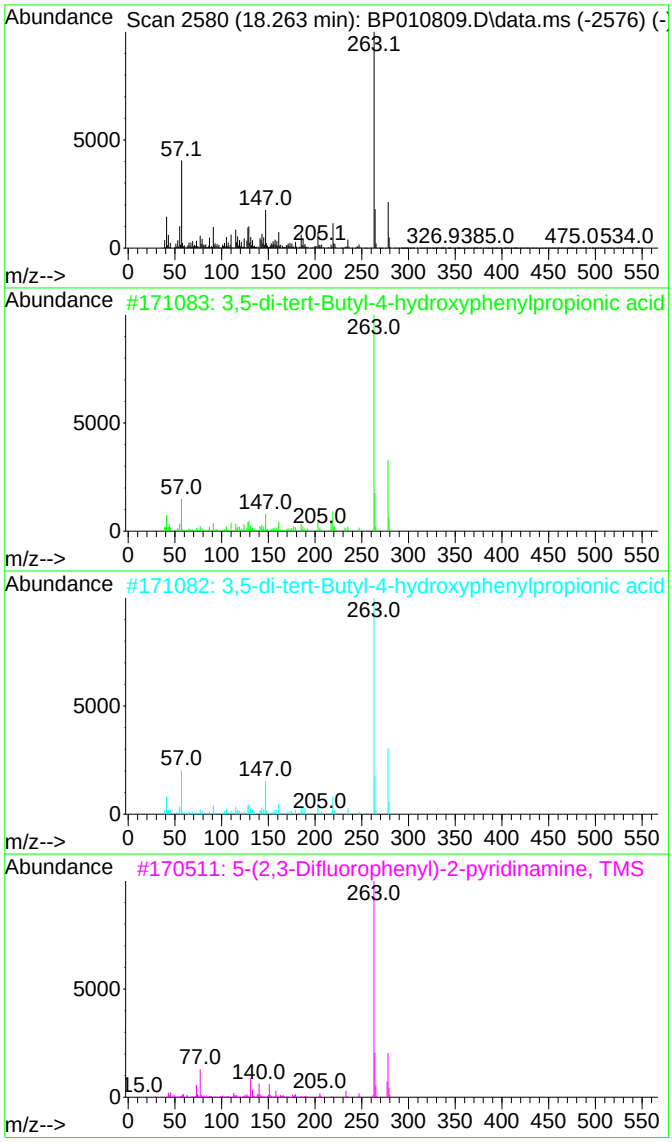
Quant Title : SVOA CALIBRATION

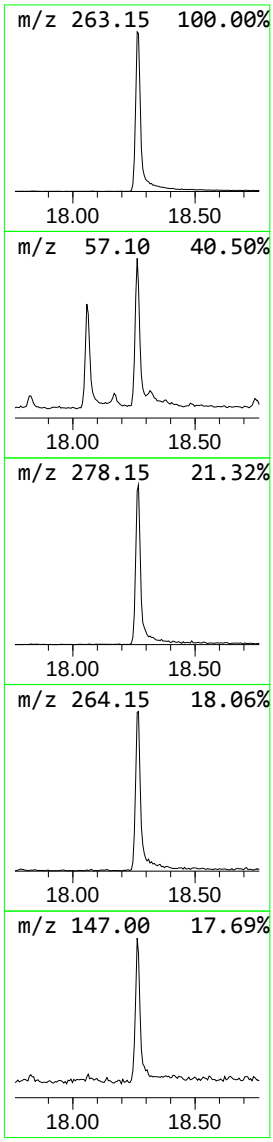
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.263	18.60 ng/ul	1832690	Phenanthrene-d10	17.139

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	93
3	5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	70
4	6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	64
5	Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	62





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ClientSampleId :
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Quant Title : SVOA CALIBRATION

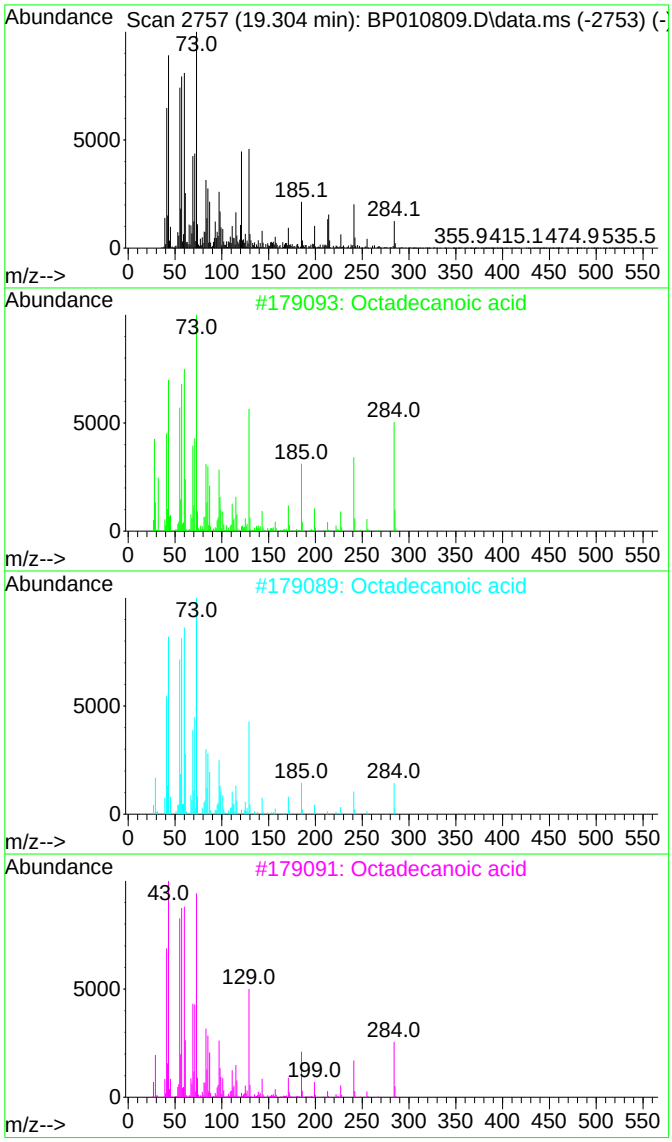
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

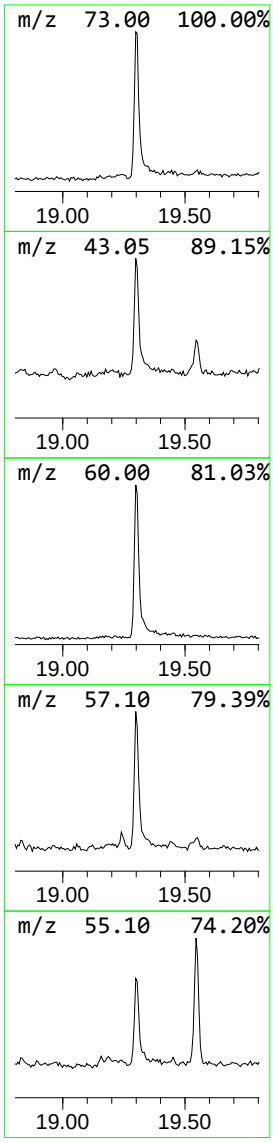
****

Peak Number 30 Octadecanoic acid Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.304	10.25 ng/ul	947517	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	93
5	Octadecanoic acid	284	C18H36O2	000057-11-4	93





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Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

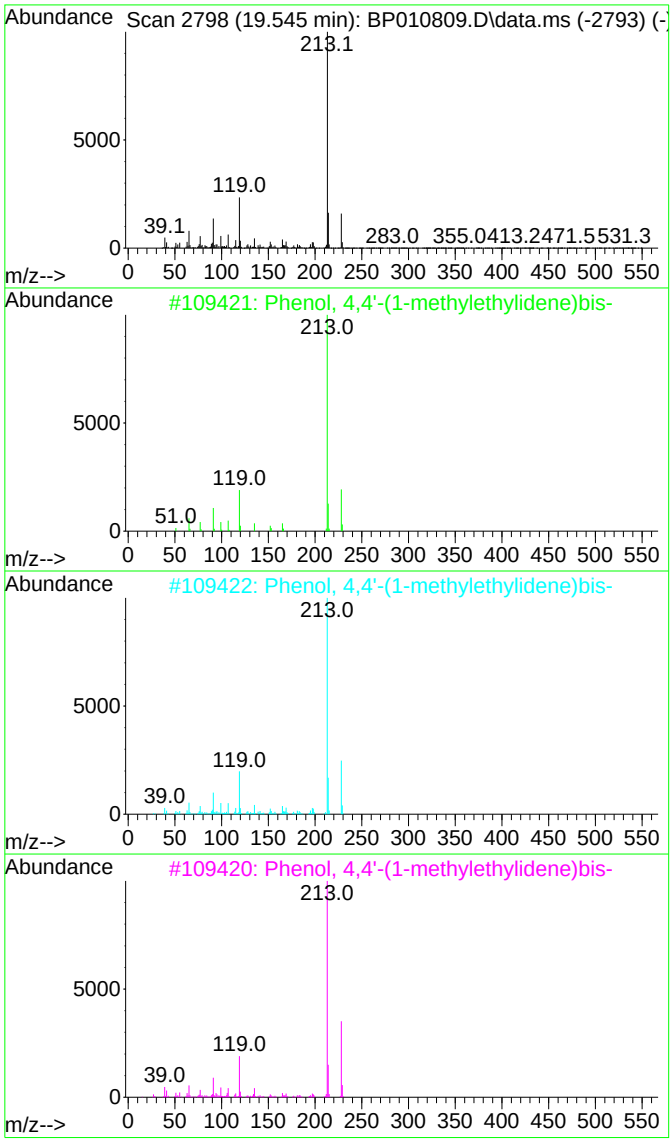
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

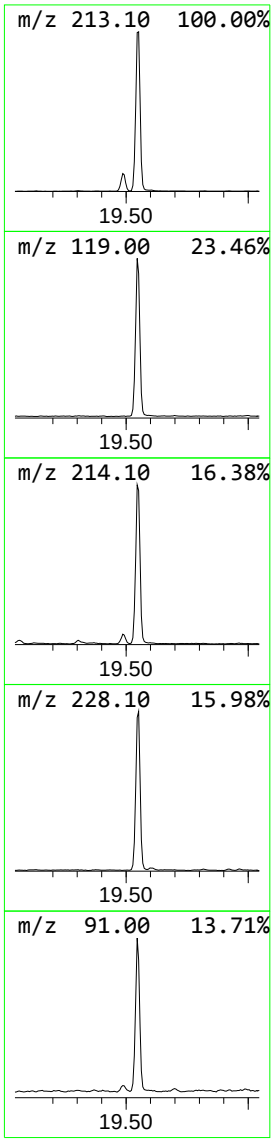
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.545	91.26 ng/ul	8432500	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	94
4	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	90
5	4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

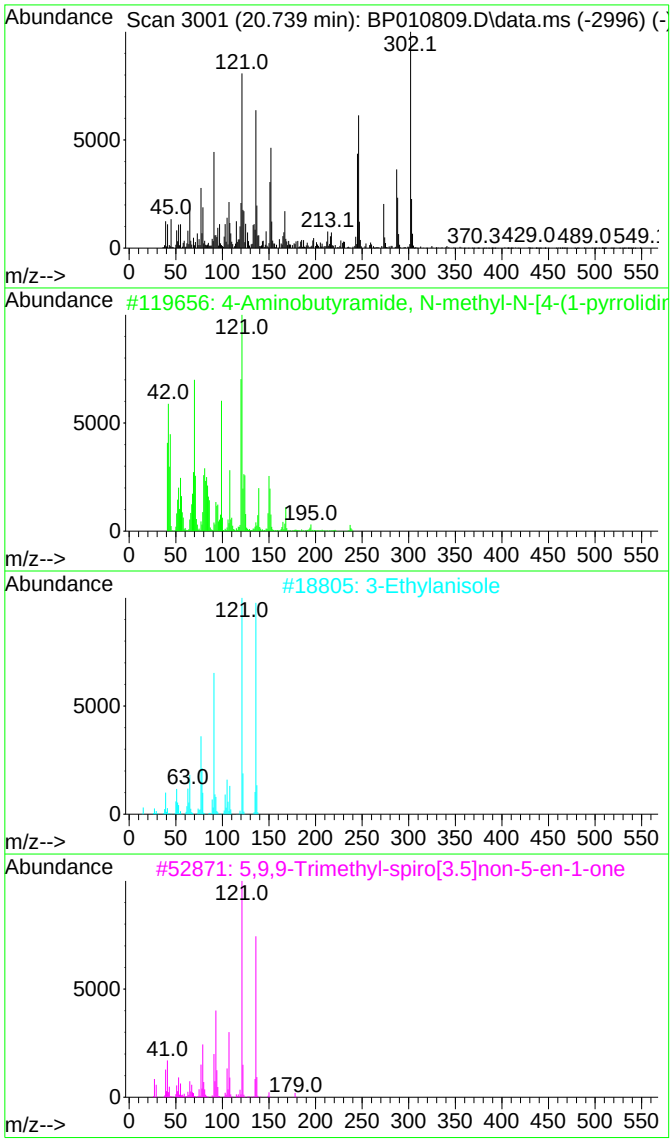
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

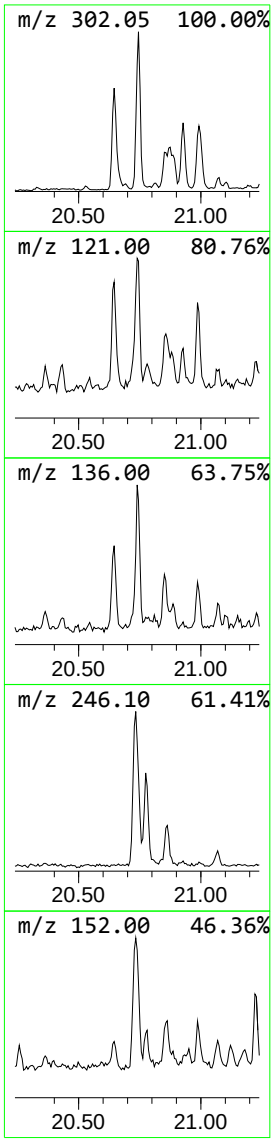
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 4-Aminobutyramide, N-methyl... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.739	8.63 ng/ul	797475	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Aminobutyramide, N-methyl-N-[4...	237	C13H23N3O	124045-56-3	53
2	3-Ethylanisole	136	C9H12O	010568-38-4	48
3	5,9,9-Trimethyl-spiro[3.5]non-5-...	178	C12H18O	1000185-13-4	20
4	trimethyl((1-methyl-4-(propan-2-...	226	C13H26OSi	1000440-50-8	20
5	3-Ethylanisole	136	C9H12O	010568-38-4	18





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

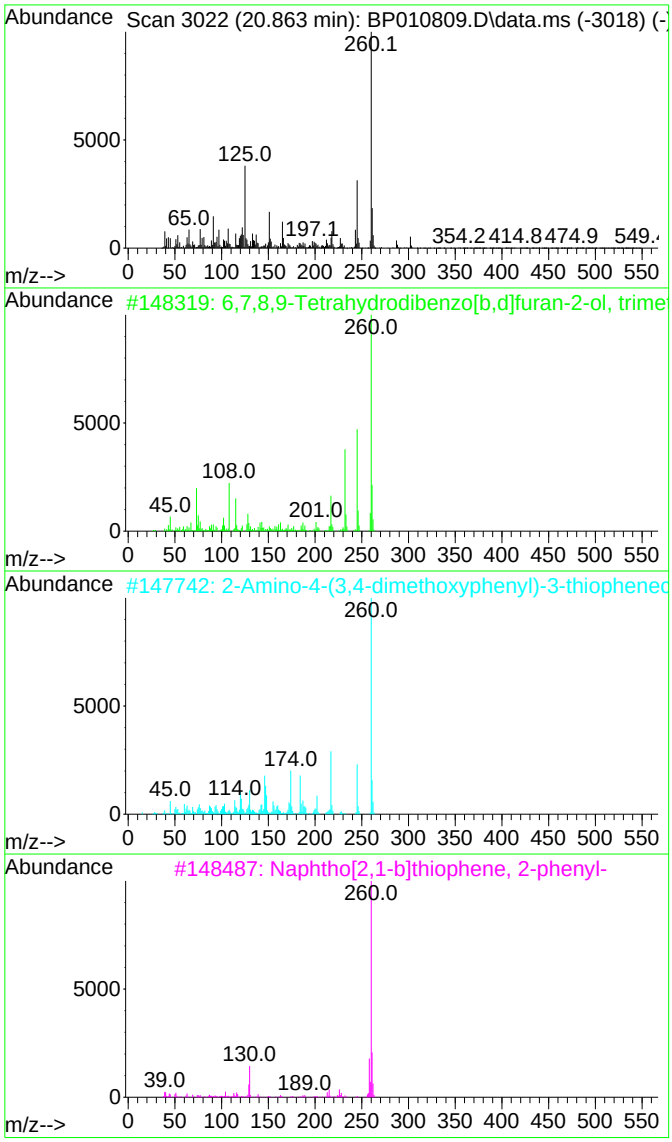
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

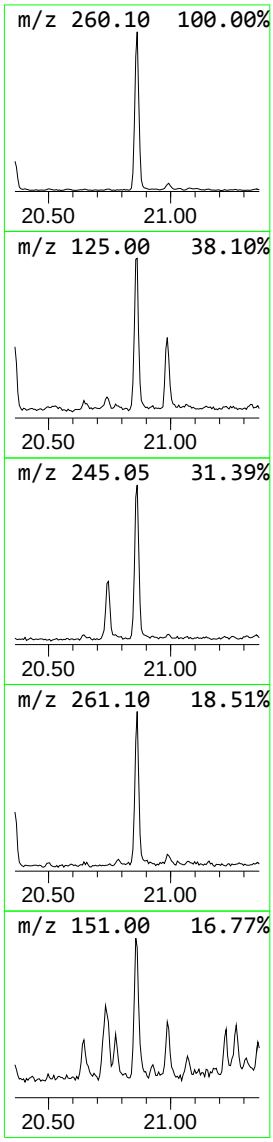
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 35 6,7,8,9-Tetrahydrodibenzo[b... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.863	17.41 ng/ul	1608770	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6,7,8,9-Tetrahydrodibenzo[b,d]fu...	260	C15H20O2Si	1000471-87-6	50
2	2-Amino-4-(3,4-dimethoxyphenyl)-...	260	C13H12N2O2S	1000475-04-7	50
3	Naphtho[2,1-b]thiophene, 2-phenyl-	260	C18H12S	016587-38-5	43
4	1,4-Benzenediamine, N,N'-diphenyl-	260	C18H16N2	000074-31-7	43
5	4-Phenyldibenzothiophene	260	C18H12S	1000314-39-7	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

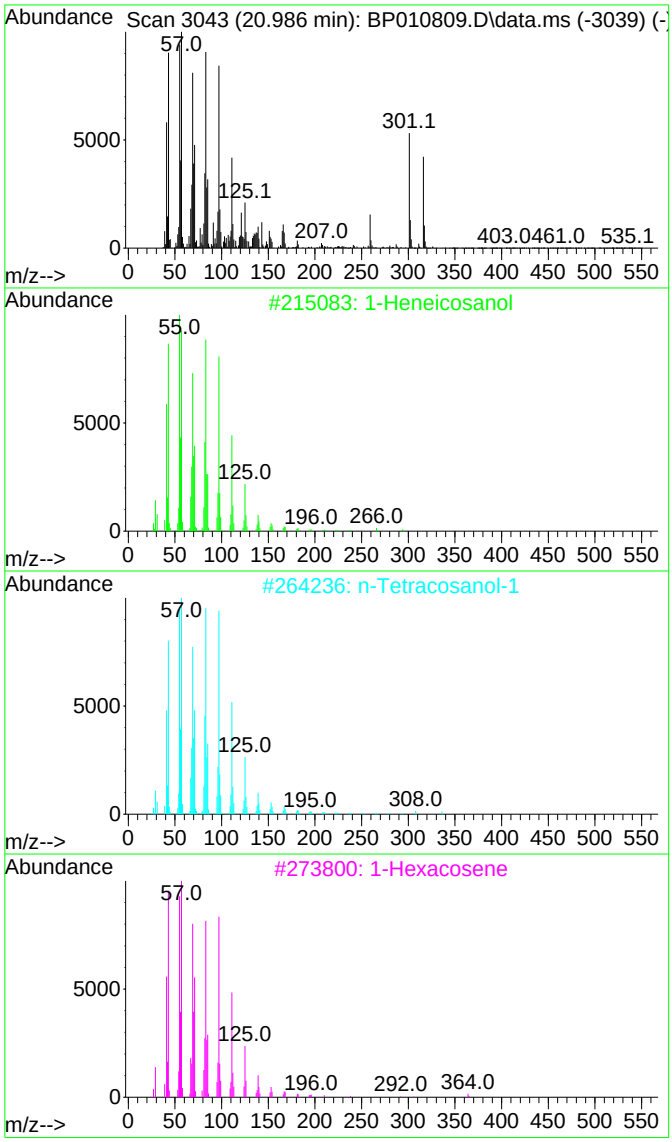
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

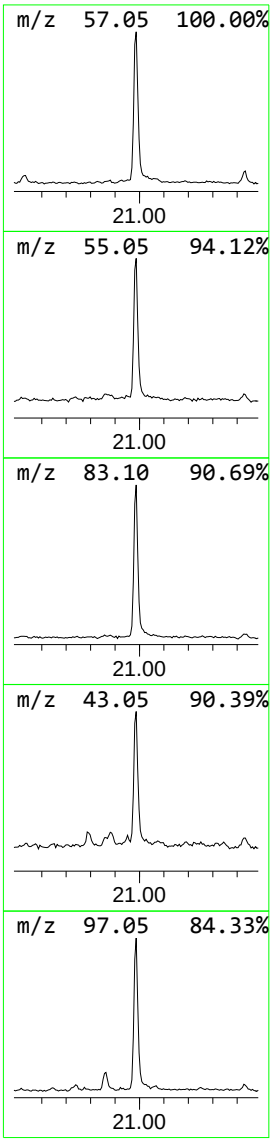
****

Peak Number 36 1-Heneicosanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	26.32 ng/ul	2432320	Chrysene-d12	21.251

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	89
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	89
3	1-Hexacosene	364	C26H52	018835-33-1	89
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	89
5	1-Tetracosene	336	C24H48	010192-32-2	89





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062422\
Data File : BP010809.D
Acq On : 25 Jun 2022 00:03
Operator : CG/JU
Sample : N3374-01
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.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-Methyl-1-hexanol	6.234	32.5	ng/ul	1713100	1	7.722	1053100	20.0
1-Hexanol, 2-et...	7.805	99.2	ng/ul	5222940	1	7.722	1053100	20.0
unknown-01	7.946	12.6	ng/ul	661702	1	7.722	1053100	20.0
Hexanoic acid, ...	9.099	18.9	ng/ul	995214	1	7.722	1053100	20.0
Phenol, p-tert-...	11.963	21.2	ng/ul	230514000	2	10.522	217721000	20.0
n-Decanoic acid	12.734	25.4	ng/ul	2326440	3	14.381	1833420	20.0
Propofol	12.775	52.3	ng/ul	4798080	3	14.381	1833420	20.0
(DEL) Alkane: S...	12.945	5.8	ng/ul	527525	3	14.381	1833420	20.0
Phenol, 2,4-bis...	13.257	1285.9	ng/ul	117881000	3	14.381	1833420	20.0
Vanillin	13.345	21.2	ng/ul	1941260	3	14.381	1833420	20.0
Phenol, 2,5-bis...	13.604	442.3	ng/ul	40542700	3	14.381	1833420	20.0
Phenol, 3,5-bis...	13.645	153.1	ng/ul	14031400	3	14.381	1833420	20.0
4-tert-Butyl-2,...	13.710	47.6	ng/ul	4363520	3	14.381	1833420	20.0
Benzofuran, 2,3...	14.151	19.1	ng/ul	1748990	3	14.381	1833420	20.0
Apocynin	14.245	17.7	ng/ul	1620590	3	14.381	1833420	20.0
1,3-Dioxolane, ...	14.881	9.5	ng/ul	874570	3	14.381	1833420	20.0
Hexathiane	14.928	9.4	ng/ul	859679	3	14.381	1833420	20.0
Phenol, 2,4,6-t...	15.063	44.7	ng/ul	4099650	3	14.381	1833420	20.0
unknown-02	15.257	12.5	ng/ul	1143630	3	14.381	1833420	20.0
Benzene, 1-(1,1...	15.551	14.1	ng/ul	1293370	3	14.381	1833420	20.0
Benzaldehyde, 4...	15.834	11.1	ng/ul	1096880	4	17.139	1970620	20.0
2(3H)-Benzothia...	16.075	16.6	ng/ul	1637260	4	17.139	1970620	20.0
Ethanone, 1-(4-...	16.439	13.4	ng/ul	1319920	4	17.139	1970620	20.0
unknown-03	16.898	13.5	ng/ul	1326500	4	17.139	1970620	20.0
n-Hexadecanoic ...	18.057	15.7	ng/ul	1548920	4	17.139	1970620	20.0
3,5-di-tert-But...	18.263	18.6	ng/ul	1832690	4	17.139	1970620	20.0
Octadecanoic acid	19.304	10.3	ng/ul	947517	5	21.251	1848020	20.0
Phenol, 4,4'-(1...	19.545	91.3	ng/ul	8432500	5	21.251	1848020	20.0
4-Aminobutyrami...	20.739	8.6	ng/ul	797475	5	21.251	1848020	20.0
6,7,8,9-Tetrahy...	20.863	17.4	ng/ul	1608770	5	21.251	1848020	20.0
1-Heneicosanol	20.986	26.3	ng/ul	2432320	5	21.251	1848020	20.0