LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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: BP010893.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	39	rVB2	530962	700182	0.25%	0.058%
2	3.646	89	95	102	rBV2	305981	1158090	0.42%	0.096%
3	6.228	528	534	545	rVV	1316284	2510852	0.91%	0.208%
4	6.928	635	653	663	rVV	18885792	34329643	12.45%	2.840%
5	7.058	669	675	685	rVB	1593465	2374422	0.86%	0.196%
6	7.246	701	707	714	rVB	2056047	3265321	1.18%	0.270%
7	7.711	777	786	793	rBV	1881168	3105535	1.13%	0.257%
8	7.799	793	801	803	rBV	1823497	3939335	1.43%	0.326%
9	7.958	819	828	843	rVB	590670	1422947	0.52%	0.118%
10	8.422	894	907	913	rBV	2133639	4273587	1.55%	0.353%
11	8.493	913	919	926	rVV	1688741	3255288	1.18%	0.269%
12	8.558	926	930	944	rVB4	325681	637285	0.23%	0.053%
13	8.828	970	976	978	rBV3	350751	615077	0.22%	0.051%
14	8.869	978	983	989	rVV	1657230	2876496	1.04%	0.238%
15	9.016	999	1008	1014	rVB8	257703	824288	0.30%	0.068%
16	9.122	1015	1026	1033	rBV2	651809	1721036	0.62%	0.142%
17	9.222	1038	1043	1050	rBV5	493441	1173890	0.43%	0.097%
18	9.340	1058	1063	1072	rBV	579205	1069785	0.39%	0.088%
19	9.434	1072	1079	1084	rBV6	605920	1619887	0.59%	0.134%
20	9.505	1084	1091	1098	rVB2	980343	2259581	0.82%	0.187%
21	9.587	1098	1105	1112	rVB3	2118102	4330864	1.57%	0.358%
22	9.699	1112	1124	1126	rBV	5918581	11025690	4.00%	0.912%
23	9.728	1127	1129	1138	rVB2	6178773	9019095	3.27%	0.746%
24	9.993	1158	1174	1178	rBV	6584974	20297853	7.36%	1.679%
25	10.034	1178	1181	1188	rVB	3901371	5937452	2.15%	0.491%
26	10.134	1189	1198	1199	rBV	2951920	4995812	1.81%	0.413%
27	10.246	1212	1217	1221	rBV	1206854	2237031	0.81%	0.185%
28	10.369	1228	1238	1243	rBV7	1560017	6058542	2.20%	0.501%
29	10.522	1243	1264	1268	rVV2	50695659	271410576	98.44%	22.450%
30	10.558	1268	1270	1280	rVB3	2976264	6579777	2.39%	0.544%
31	10.705	1287	1295	1302	rBV3	2775965	6841279	2.48%	0.566%
32	10.769	1302	1306	1307	rBV2	858149	1031318	0.37%	0.085%
33	10.805	1308	1312	1319	rVB2	3199652	5755033	2.09%	0.476%
34	10.905	1319	1329	1334	rBV2	36515953	80383713	29.15%	6.649%
35	10.981	1340	1342	1349	rVB3	1439729	2290647	0.83%	0.189%
36	11.069	1350	1357	1361	rBV	1259494	2119249	0.77%	0.175%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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	11.140	1361	1369	1374	rVB2	2417178	4690773	1.70%	0.388%
38	11.275	1386	1392	1396	rBV4	546541	1248750	0.45%	0.103%
39	11.346	1397	1404	1409	rVB5	1261549	2603108	0.94%	0.215%
40	11.440	1409	1420	1423	rBV7	1415946	4687813	1.70%	0.388%
41	11.540	1426	1437	1441	rVB4	4641264	10727595	3.89%	0.887%
42	11.587	1441	1445	1449	rVB2	2646107	3702510	1.34%	0.306%
43	11.622	1449	1451	1455	rVB2	801284	916844	0.33%	0.076%
44	11.693	1455	1463	1468	rBV5	3941406	10905211	3.96%	0.902%
45	11.769	1471	1476	1478	rBV3	2137440	3604586	1.31%	0.298%
46	11.963	1489	1509	1513	rVB3	48903874	275720778	100.00%	22.806%
47	12.034	1517	1521	1524	rBV2	1564587	2244068	0.81%	0.186%
48	12.075	1524	1528	1530	rBV	1518396	2268383	0.82%	0.188%
49	12.146	1535	1540	1546	rVB	2789896	5189168	1.88%	0.429%
50	12.257	1546	1559	1567	rBV6	3478155	12231340	4.44%	1.012%
51	12.516	1595	1603	1608	rBV6	339405	816372	0.30%	0.068%
52	12.569	1608	1612	1620	rBV6	352409	697827	0.25%	0.058%
53	12.763	1641	1645	1658	rVB	4363121	7317479	2.65%	0.605%
54	12.940	1672	1675	1683	rVB7	286520	676715	0.25%	0.056%
55	13.251	1716	1728	1732	rBV3	53112078	145193059	52.66%	12.010%
56	13.340	1739	1743	1750	rVB	2110337	2831066	1.03%	0.234%
57	13.499	1766	1770	1777	rVB	462177	703074	0.25%	0.058%
58	13.599	1778	1787	1790	rBV	34170594	54562374	19.79%	4.513%
59	13.634	1790	1793	1800	rVV	14142280	19428067	7.05%	1.607%
60	13.704	1800	1805	1811	rVB	4296161	5881373	2.13%	0.486%
61	13.799	1817	1821	1825	rVV	2076069	2570861	0.93%	0.213%
62	13.840	1825	1828	1836	rVB	983161	1464768	0.53%	0.121%
63	14.063	1859	1866	1874	rVB	3090068	5002049	1.81%	0.414%
64	14.140	1874	1879	1883	rBV	2105502	2719804	0.99%	0.225%
65	14.234	1891	1895	1899	rBV2	1479038	2046212	0.74%	0.169%
66	14.310	1905	1908	1913	rVB2	617509	810420	0.29%	0.067%
67	14.369	1913	1918	1923	rBV	3027708	4607032	1.67%	0.381%
68	14.422	1923	1927	1936	rVB	1967167	2654428	0.96%	0.220%
69	14.581	1950	1954	1956	rBV2	459911	624256	0.23%	0.052%
70	14.616	1956	1960	1965	rVB	3368517	4397264	1.59%	0.364%
71	14.869	1999	2003	2007	rVB	681019	865788	0.31%	0.072%
72	14.922	2007	2012	2019	rVB3	491141	968721	0.35%	0.080%
73	15.051	2029	2034	2039	rBV	4104885	5279720	1.91%	0.437%
74	15.245	2063	2067	2071	rVV	1254450	1513321	0.55%	0.125%
75	15.363	2082	2087	2095	rBV	3808902	5742327	2.08%	0.475%
76	15.487	2103	2108	2113	rBV2	1355850	2145571	0.78%	0.177%

LSC Area Percent Report

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Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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	15.545	2113	2118	2124	rVB	1211134	1689838	0.61%	0.140%
78	15.634	2127	2133	2141	rVB7	308946	675646	0.25%	0.056%
79	15.828	2162	2166	2171	rVB	1093112	1412311	0.51%	0.117%
80	16.075	2201	2208	2215	rBV	1282465	2362248	0.86%	0.195%
81	16.434	2264	2269	2281	rVB	1542936	2368413	0.86%	0.196%
82	16.892	2342	2347	2353	rVB	1173472	1793923	0.65%	0.148%
83	17.128	2382	2387	2397	rVB	3250101	4926640	1.79%	0.408%
84	17.228	2399	2404	2412	rVV	4132992	5721971	2.08%	0.473%
85	17.516	2449	2453	2456	rBV	534909	787020	0.29%	0.065%
86	18.045	2539	2543	2552	rBV	1062284	1553467	0.56%	0.128%
87	18.251	2574	2578	2585	rBV	1023879	1526428	0.55%	0.126%
88	19.286	2751	2754	2763	rBV3	434031	730270	0.26%	0.060%
89	19.481	2782	2787	2792	rBV	4830713	6341887	2.30%	0.525%
90	19.539	2792	2797	2805	rBV	9841561	11967237	4.34%	0.990%
91	20.086	2887	2890	2901	rVB6	556982	1009399	0.37%	0.083%
92	20.351	2932	2935	2939	rBV	1035230	1190126	0.43%	0.098%
93	20.633	2980	2983	2989	rVB3	441286	613965	0.22%	0.051%
94	20.728	2995	2999	3003	rVB3	875342	1115692	0.40%	0.092%
95	20.851	3016	3020	3027	rBV2	1491574	2150891	0.78%	0.178%
96	20.969	3036	3040	3048	rBV3	2541205	3433242	1.25%	0.284%
97	21.239	3082	3086	3093	rVB	4044096	5203276	1.89%	0.430%
98	21.516	3130	3133	3138	rVB	966662	1185043	0.43%	0.098%
99	23.457	3456	3463	3472	rVB2	3875468	7567516	2.74%	0.626%
100	23.610	3482	3489	3502	rVB2	2883768	5934003	2.15%	0.491%

Sum of corrected areas: 1208967785

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
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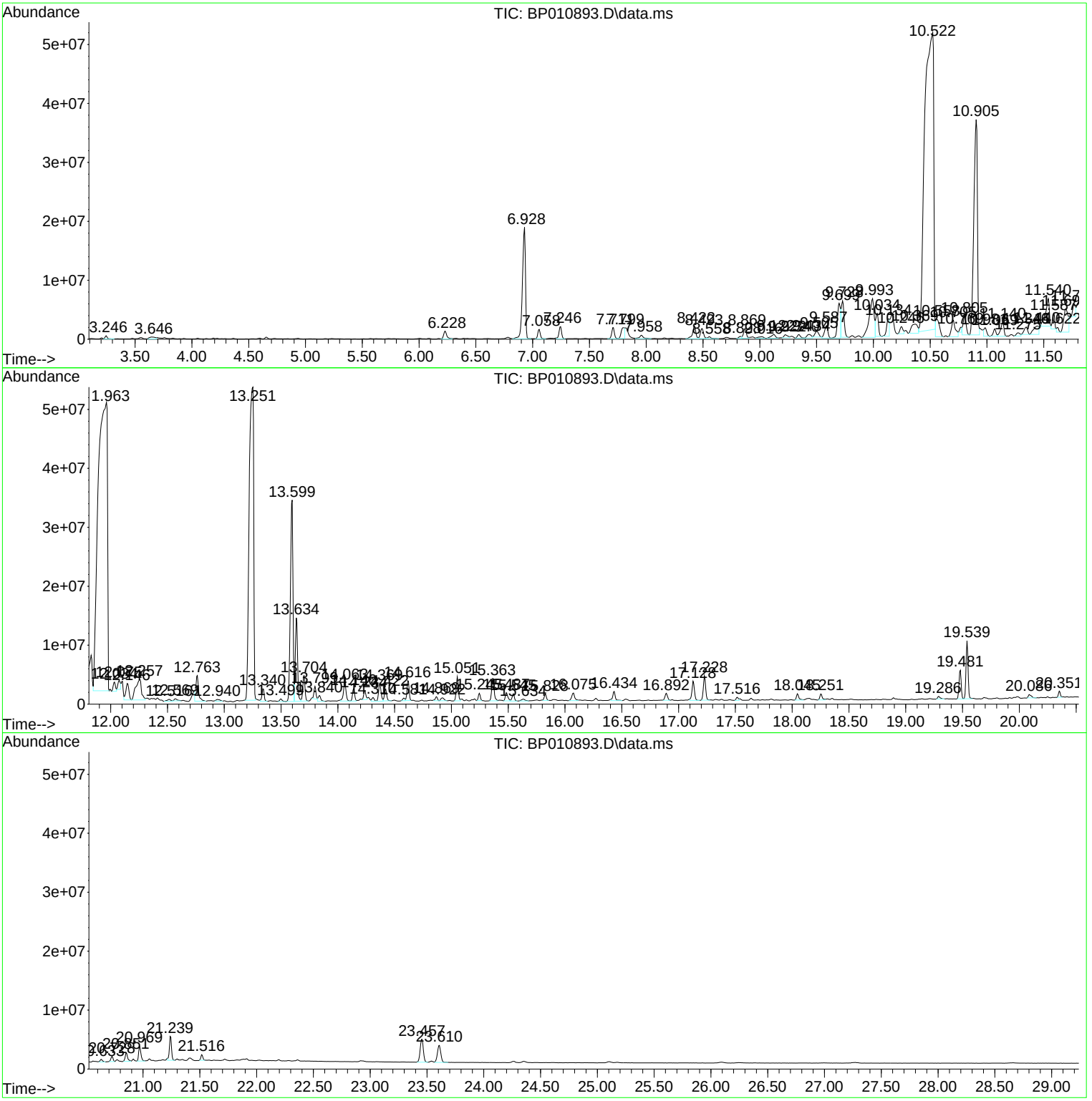
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

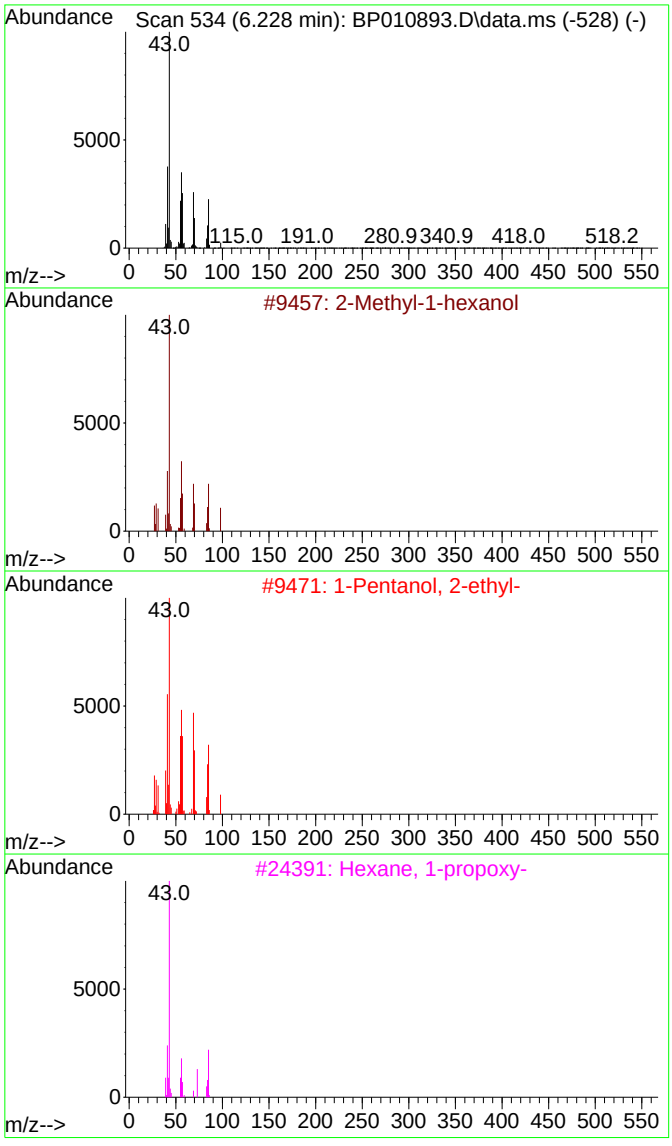
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

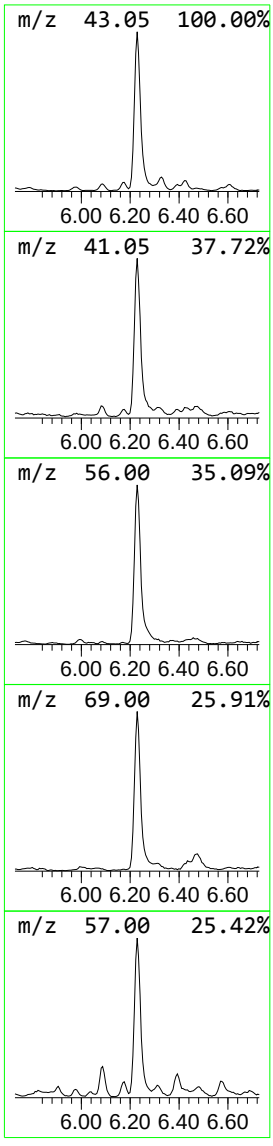
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Methyl-1-hexanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.228	16.17 ng/ul	2510850	1,4-Dichlorobenzene-d4	7.711

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-1-hexanol	116	C7H16O	1000427-67-0	78
2	1-Pentanol, 2-ethyl-	116	C7H16O	027522-11-8	59
3	Hexane, 1-propoxy-	144	C9H20O	053685-78-2	59
4	Octane	114	C8H18	000111-65-9	53
5	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

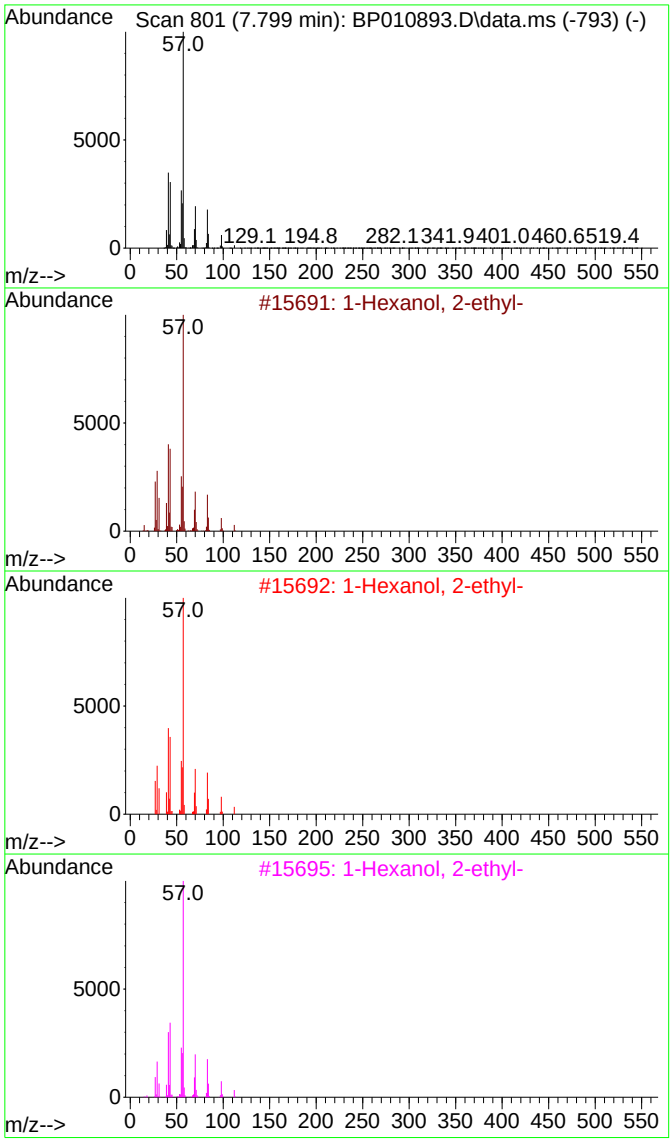
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

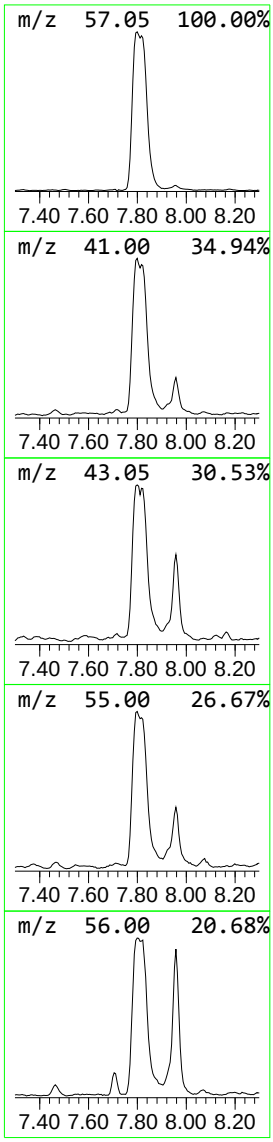
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.799	25.37 ng/ul	3939340	1,4-Dichlorobenzene-d4	7.711

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	78
4	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
5	2-Ethyl-1-hexanol, methyl ether	144	C9H20O	1000365-10-2	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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

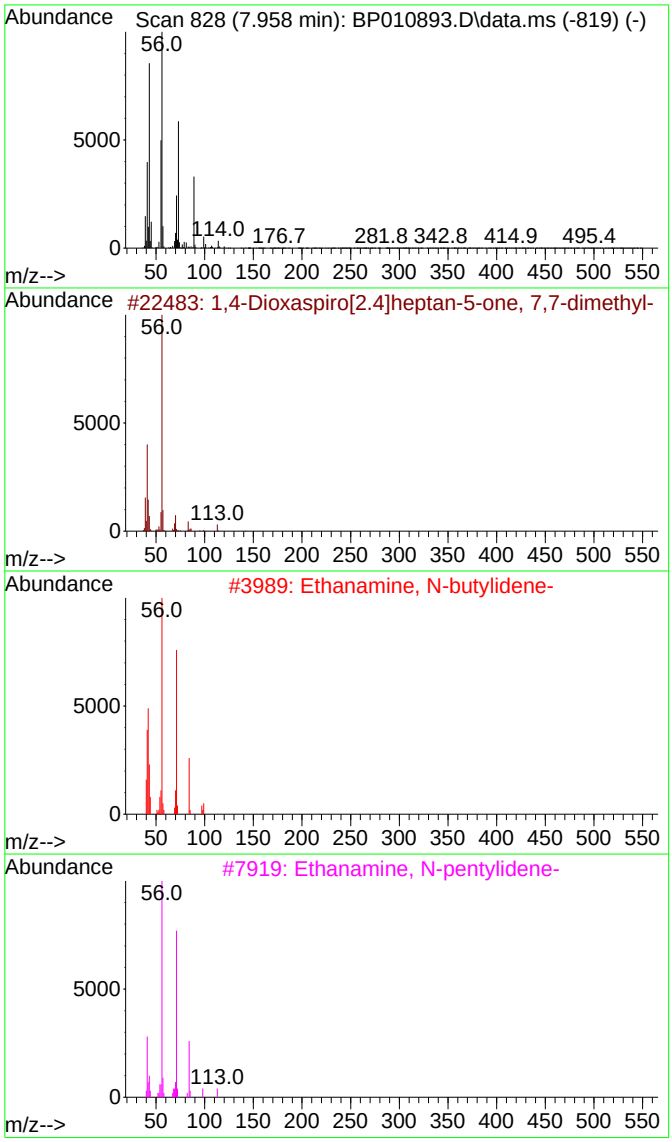
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

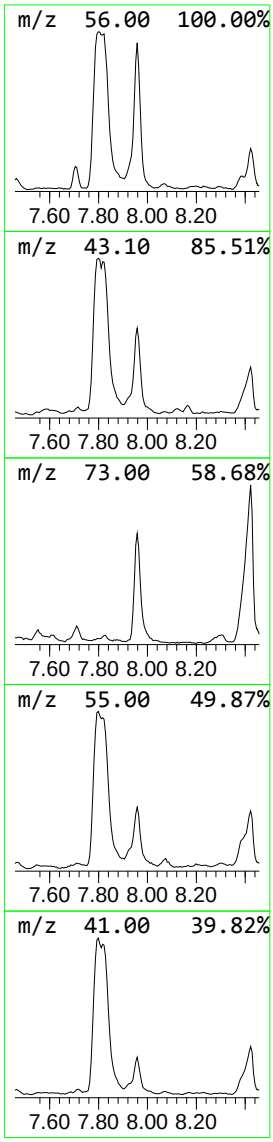
****

Peak Number 3 unknown-01 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.958	9.16 ng/ul	1422950	1,4-Dichlorobenzene-d4	7.711

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	46
2	Ethanamine, N-butylidene-	99	C6H13N	001611-12-7	43
3	Ethanamine, N-pentylidene-	113	C7H15N	010599-76-5	40
4	2(3H)-Furanone, dihydro-3,3-dime...	114	C6H10O2	003709-08-8	37
5	4-Aminocyclohexanol acetate (ester)	157	C8H15NO2	1000130-62-8	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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

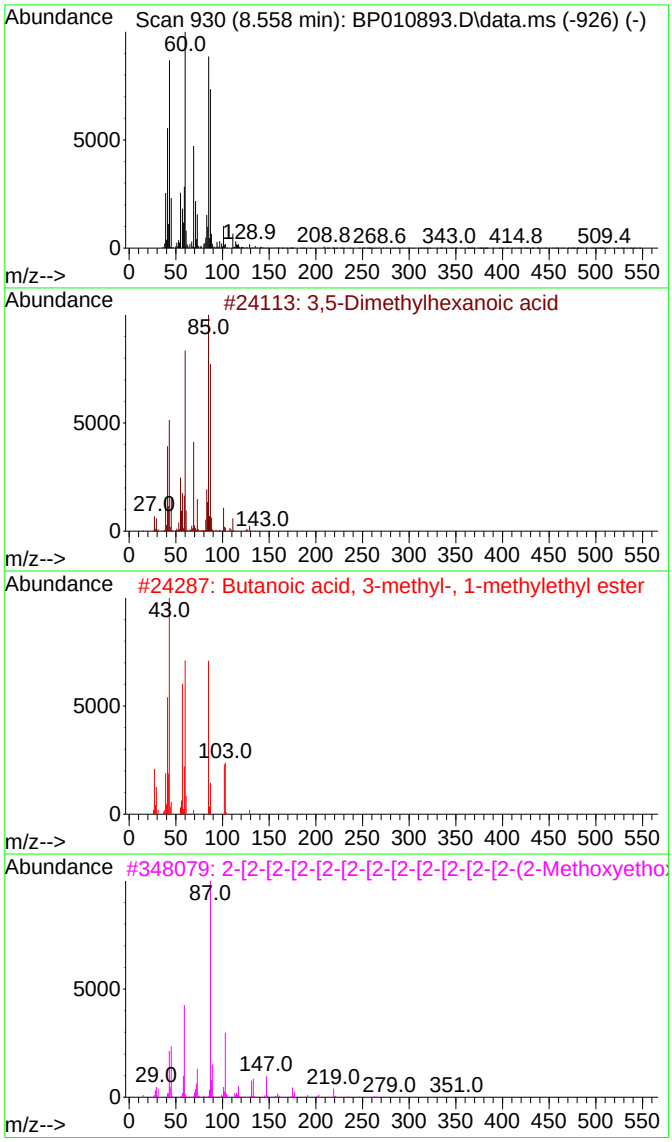
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

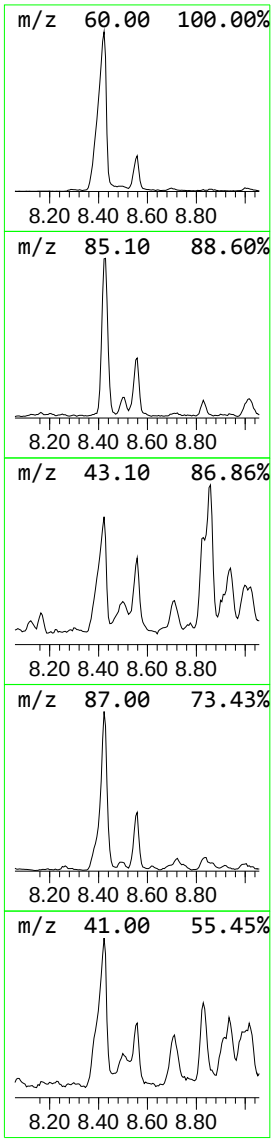
****

Peak Number 4 3,5-Dimethylhexanoic acid Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.558	4.10 ng/ul	637285	1,4-Dichlorobenzene-d4	7.711

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethylhexanoic acid	144	C8H16O2	060308-87-4	53
2	Butanoic acid, 3-methyl-, 1-meth...	144	C8H16O2	032665-23-9	43
3	2-[2-[2-[2-[2-[2-[2-[2-[2-[2-...	646	C29H58O15	1000351-92-0	27
4	Diallylmethylsilane	126	C7H14Si	002043-08-5	27
5	1,3-Dioxane, 2-methyl-	102	C5H10O2	000626-68-6	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

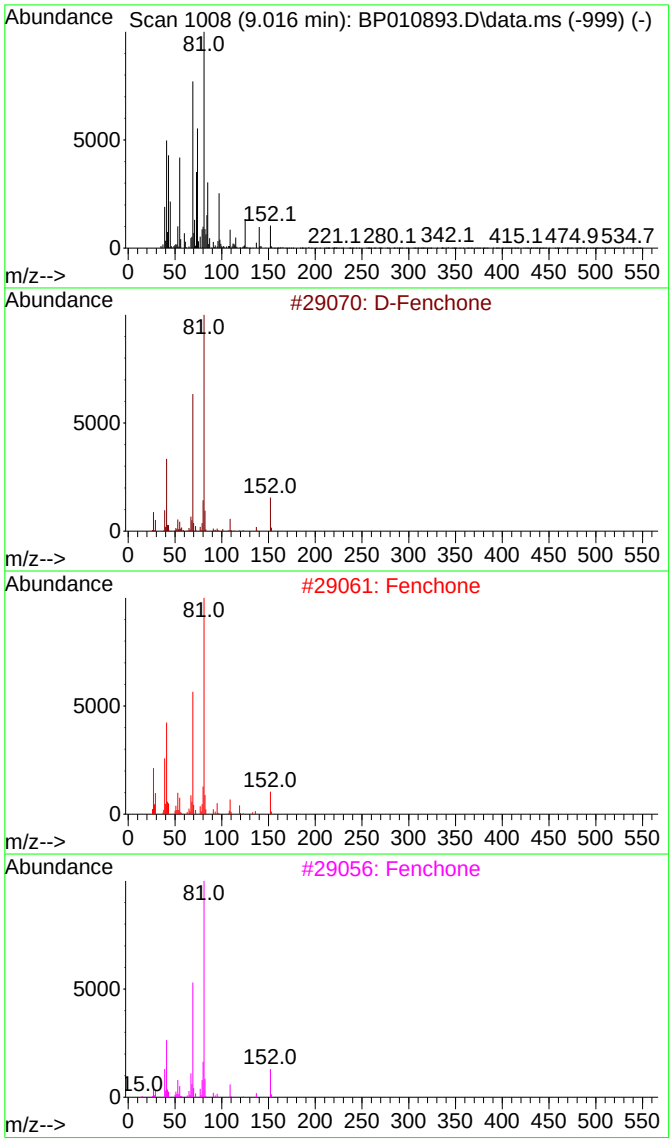
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

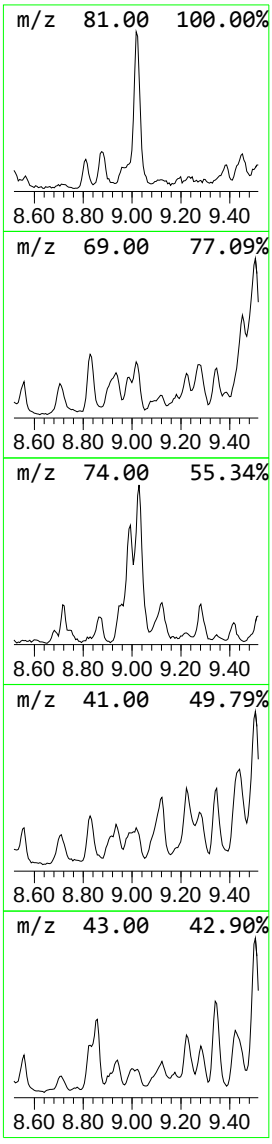
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown-02 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.016	5.31 ng/ul	824288	1,4-Dichlorobenzene-d4	7.711

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Fenchone	152	C10H16O	004695-62-9	43
2	Fenchone	152	C10H16O	001195-79-5	43
3	Fenchone	152	C10H16O	001195-79-5	32
4	3-Ethyl-4-octene	140	C10H20	019781-32-9	22
5	Cyclopentanecarboxylic acid, cyc...	168	C10H16O2	1000282-76-8	22





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Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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

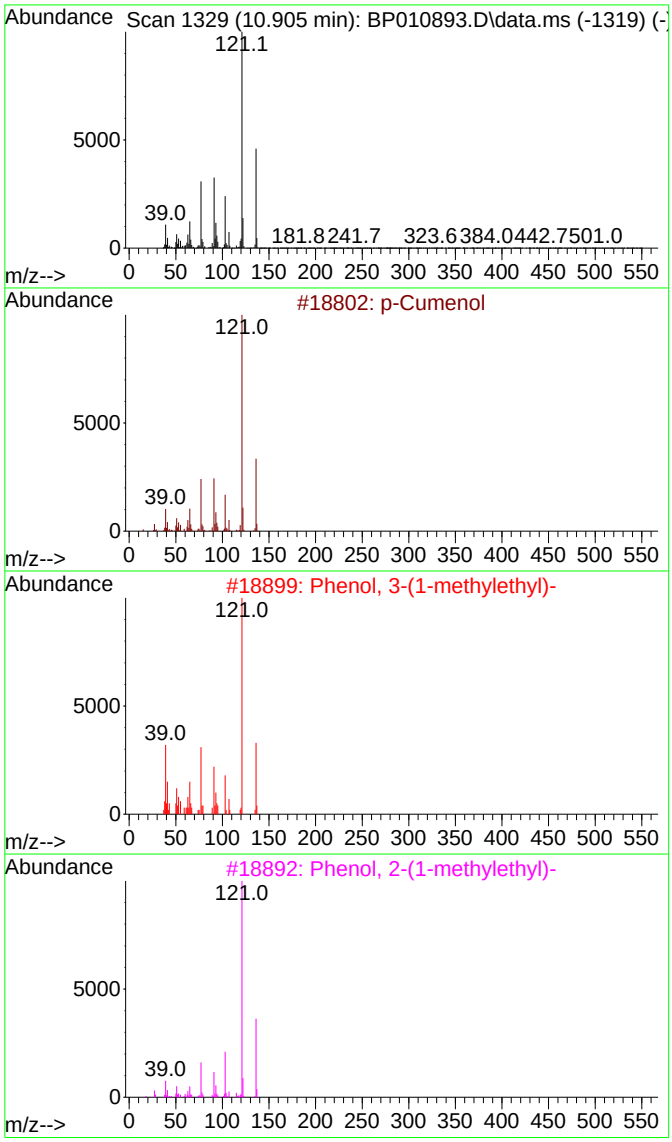
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

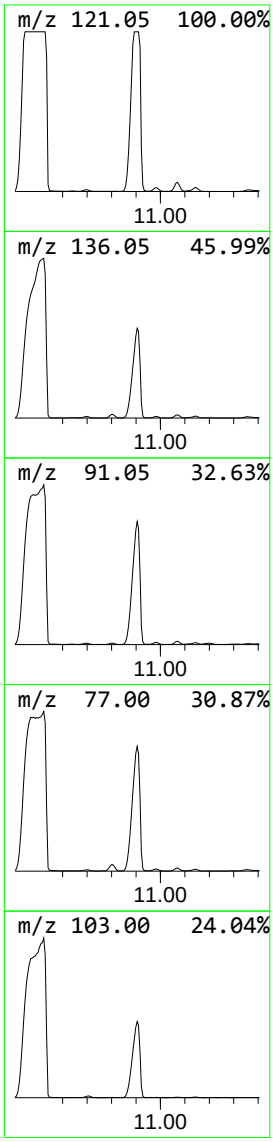
****

Peak Number 6 p-Cumenol Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.905	5.92 ng/ul	80383700	Naphthalene-d8	10.522

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cumenol	136	C9H12O	000099-89-8	96
2	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91
3	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	87
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	87





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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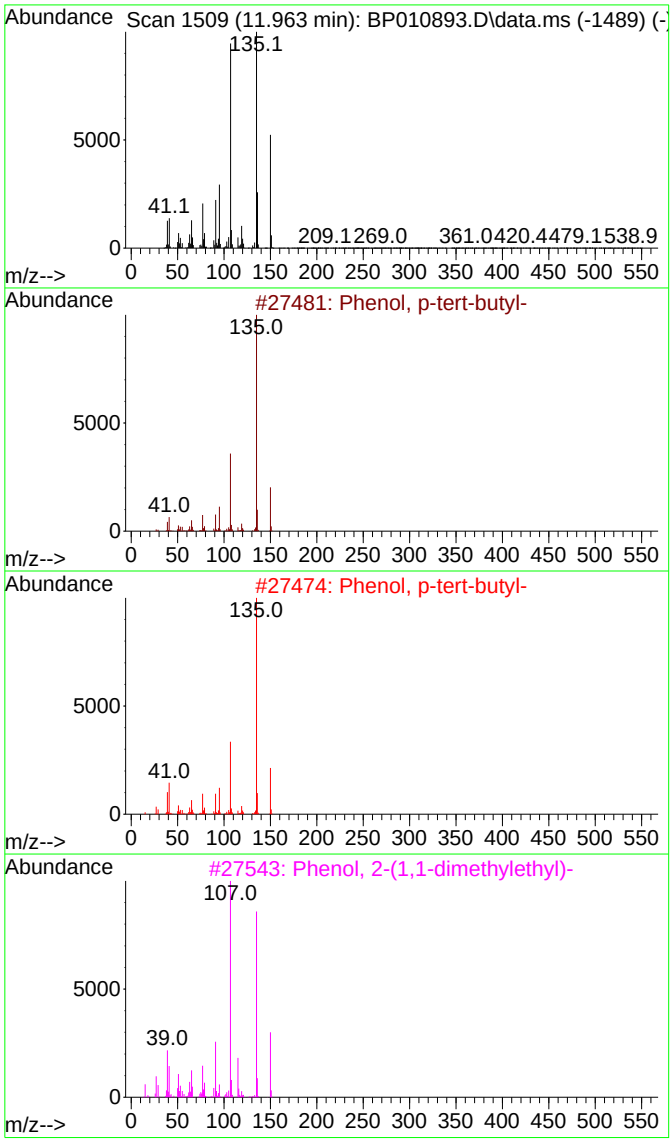
Quant Title : SVOA CALIBRATION

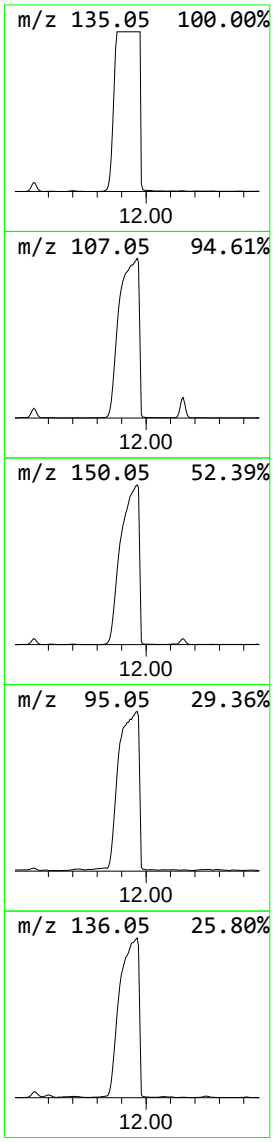
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Phenol, p-tert-butyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	20.32 ng/ul	275721000	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	74
4	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	68
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

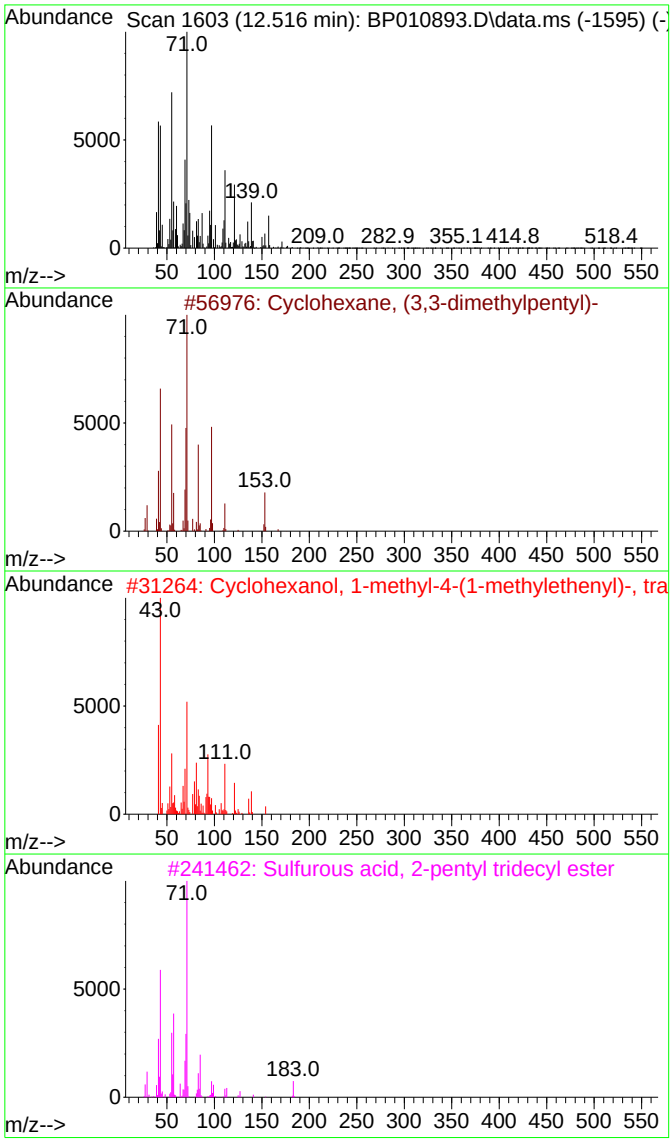
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

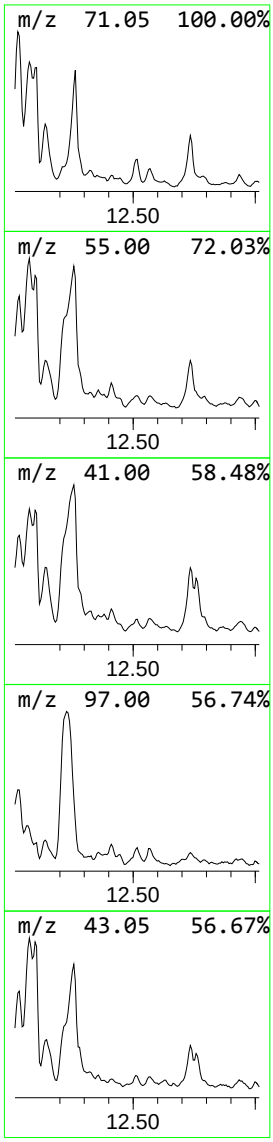
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Cyclic12.516 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.516	3.54 ng/ul	816372	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, (3,3-dimethylpentyl)-	182	C13H26	061142-22-1	43
2	Cyclohexanol, 1-methyl-4-(1-meth...	154	C10H18O	007299-40-3	38
3	Sulfurous acid, 2-pentyl tridecy...	334	C18H38O3S	1000309-16-2	27
4	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	25
5	Sulfurous acid, 2-pentyl tetrade...	348	C19H40O3S	1000309-16-3	22





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

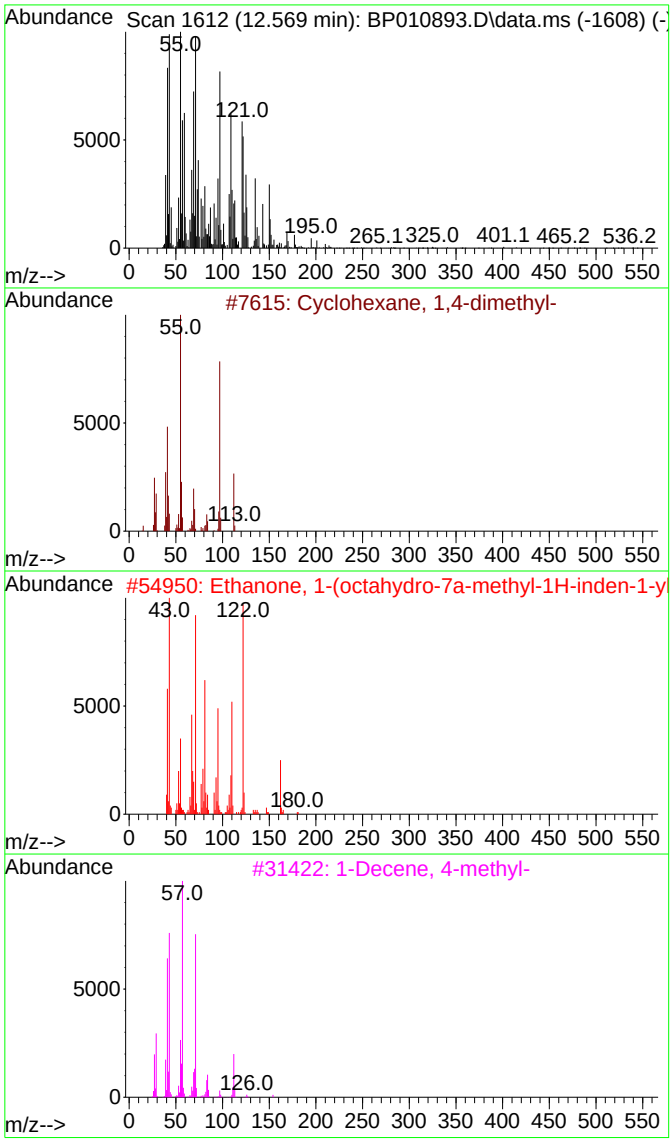
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

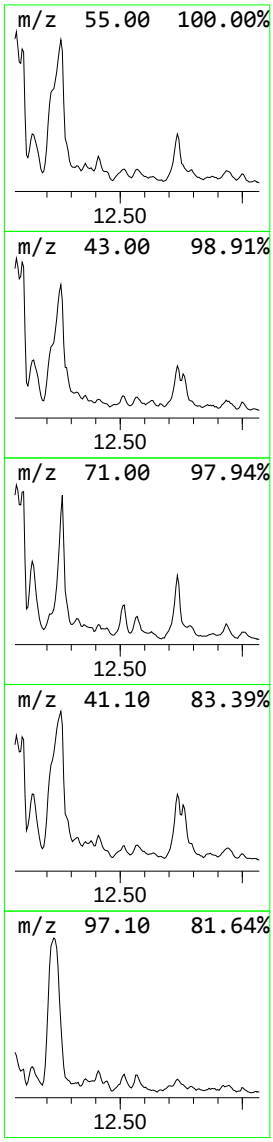
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Cyclic12.569 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.569	3.03 ng/ul	697827	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	25
2	Ethanone, 1-(octahydro-7a-methyl...	180	C12H20O	017986-96-8	25
3	1-Decene, 4-methyl-	154	C11H22	013151-29-6	15
4	Nonane, 5-butyl-	184	C13H28	017312-63-9	11
5	3-Hepten-2-one	112	C7H12O	001119-44-4	11





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

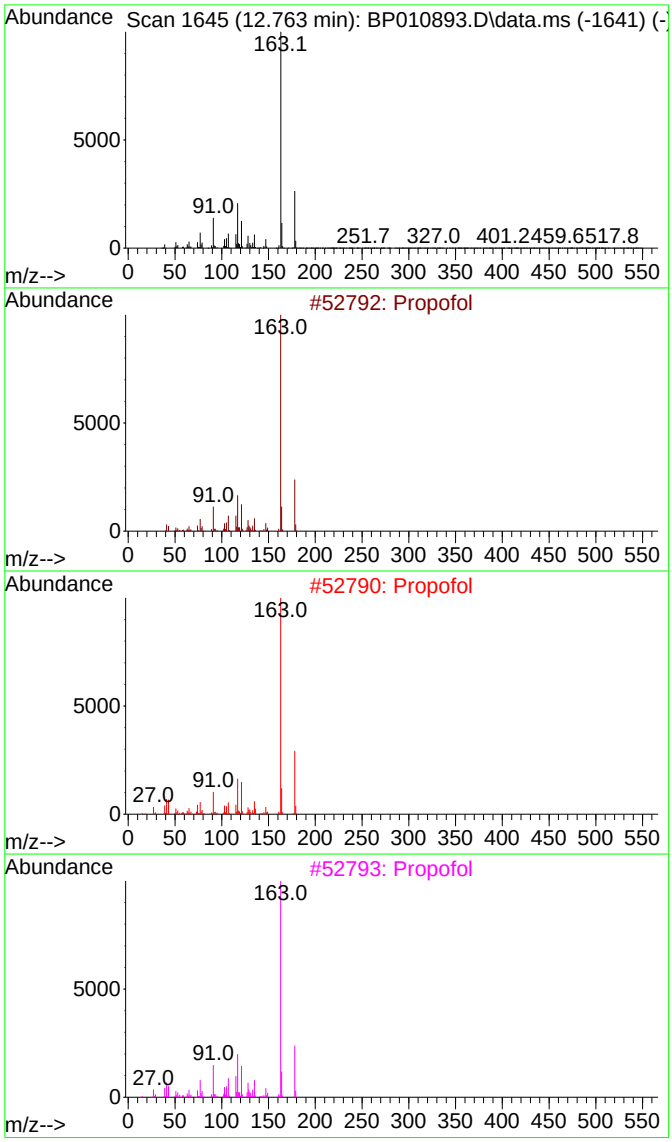
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

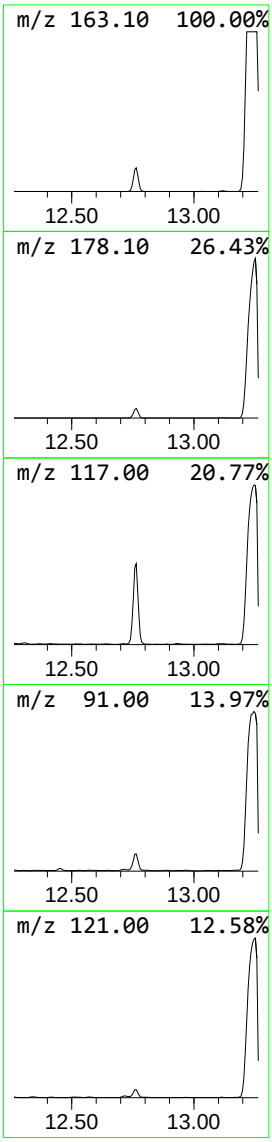
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Propofol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.763	31.77 ng/ul	7317480	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	96
2	Propofol	178	C12H18O	002078-54-8	94
3	Propofol	178	C12H18O	002078-54-8	93
4	Propofol	178	C12H18O	002078-54-8	93
5	Propofol	178	C12H18O	002078-54-8	90





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

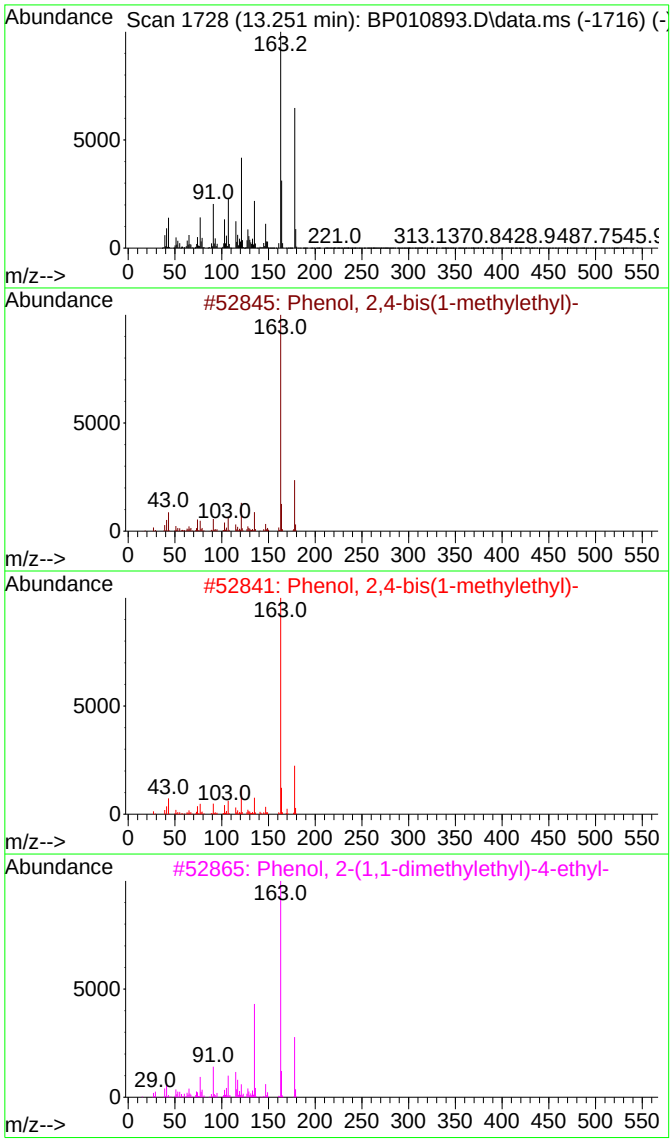
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

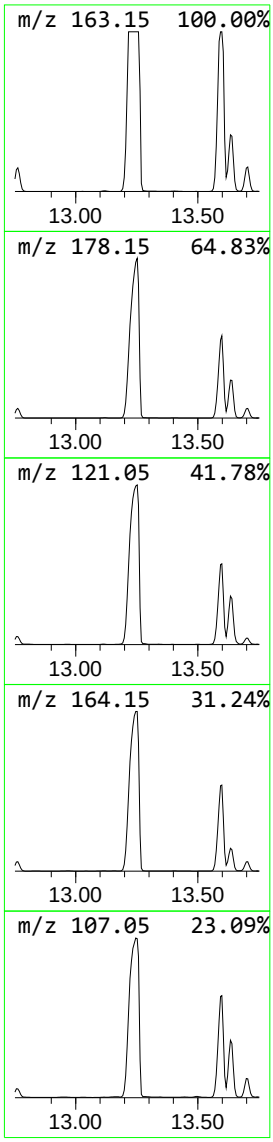
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Phenol, 2,4-bis(1-methyleth... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.252	630.31 ng/ul	145193000	Acenaphthene-d10	14.369

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	Phenol, 2-(1,1-dimethylethyl)-4-...	178	C12H18O	000096-70-8	72
4	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	70
5	Propofol	178	C12H18O	002078-54-8	55





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Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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

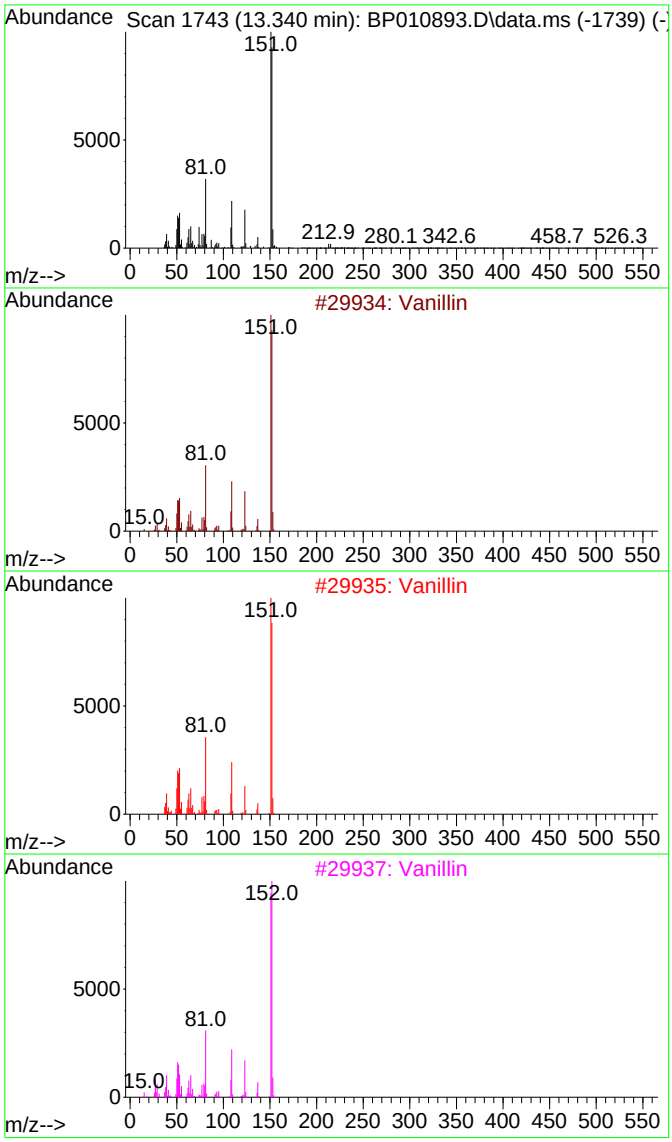
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

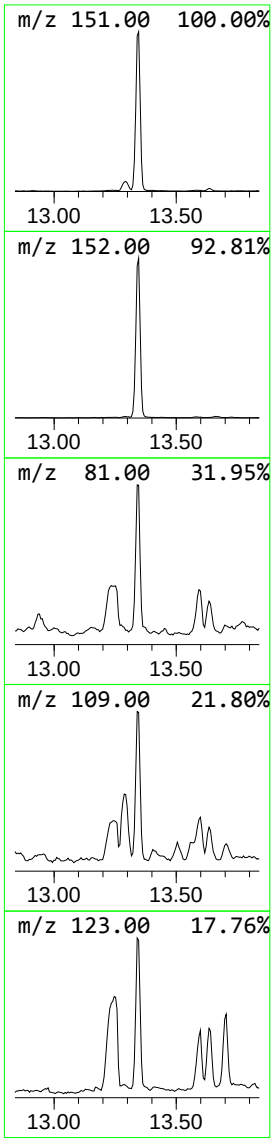
****

Peak Number 13 Vanillin Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.340	12.29 ng/ul	2831070	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	97
2	Vanillin	152	C8H8O3	000121-33-5	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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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

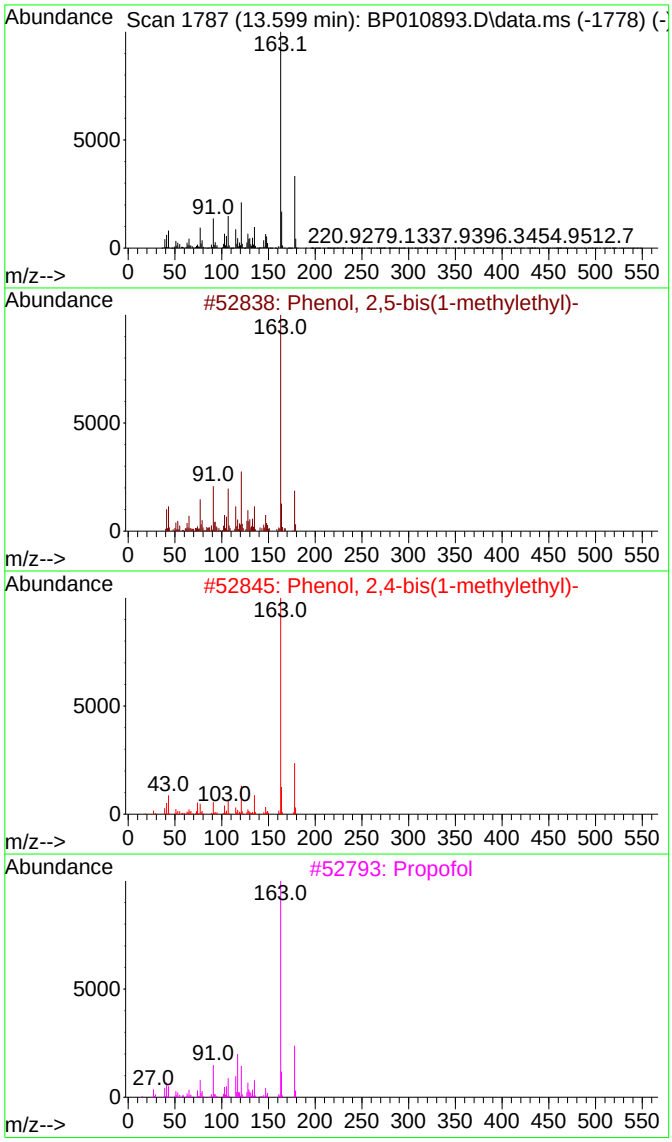
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

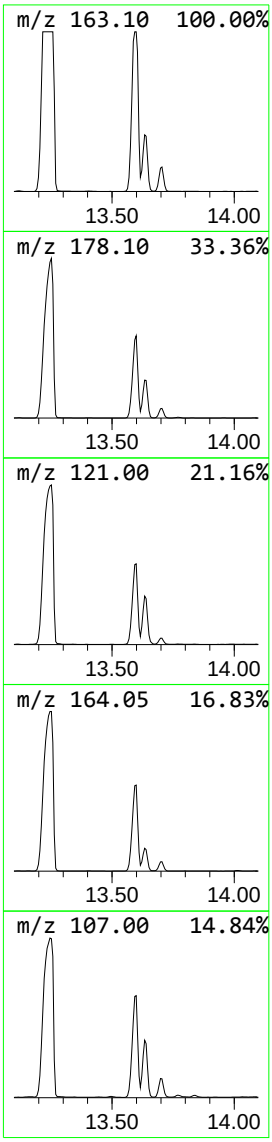
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Phenol, 2,5-bis(1-methyleth... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.599	236.87 ng/ul	54562400	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,5-bis(1-methylethyl)-	178	C12H18O	035946-91-9	94
2	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	90
3	Propofol	178	C12H18O	002078-54-8	87
4	Phenol, 2,4-bis(1-methylethyl)-	178	C12H18O	002934-05-6	87
5	Propofol	178	C12H18O	002078-54-8	87





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

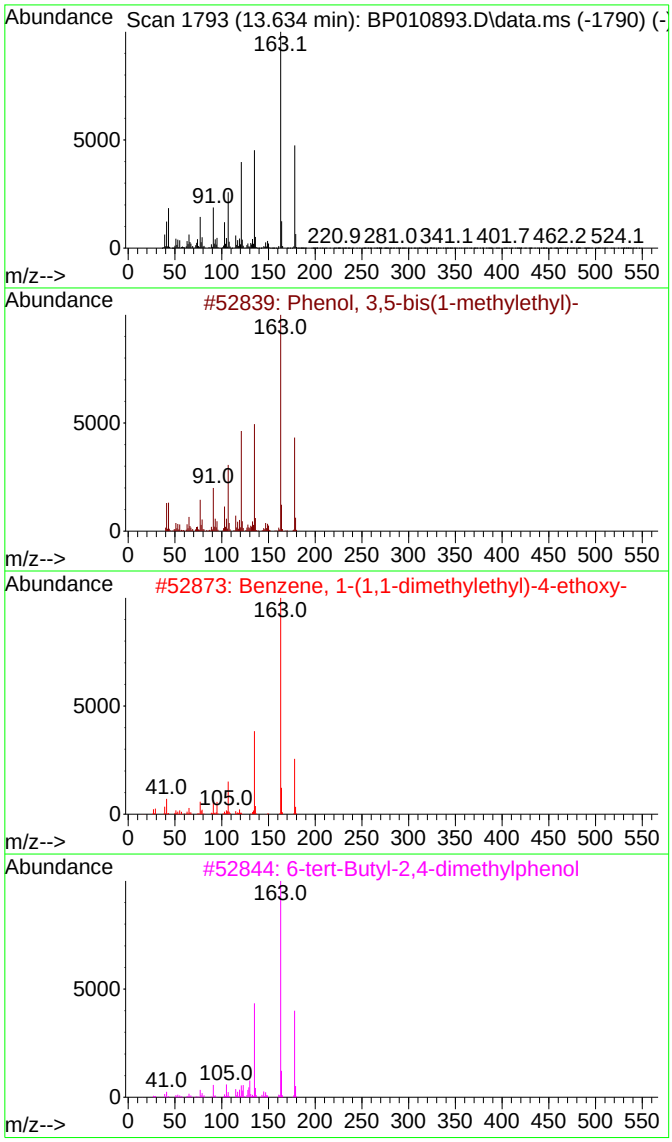
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

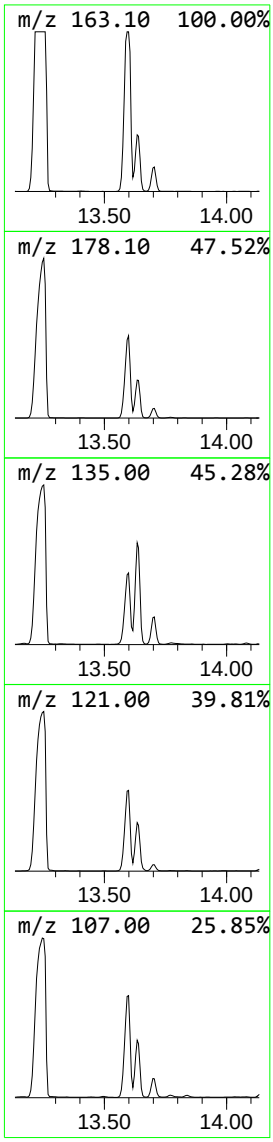
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Phenol, 3,5-bis(1-methyleth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.634	84.34 ng/ul	19428100	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-bis(1-methylethyl)-	178	C12H18O	026886-05-5	96
2	Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	90
3	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	81
4	5'-Hydroxy-2',3',4'-trimethylace...	178	C11H14O2	013667-28-2	81
5	Phenol, 2-(1,1-dimethylethyl)-4-...	178	C12H18O	000096-70-8	72





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

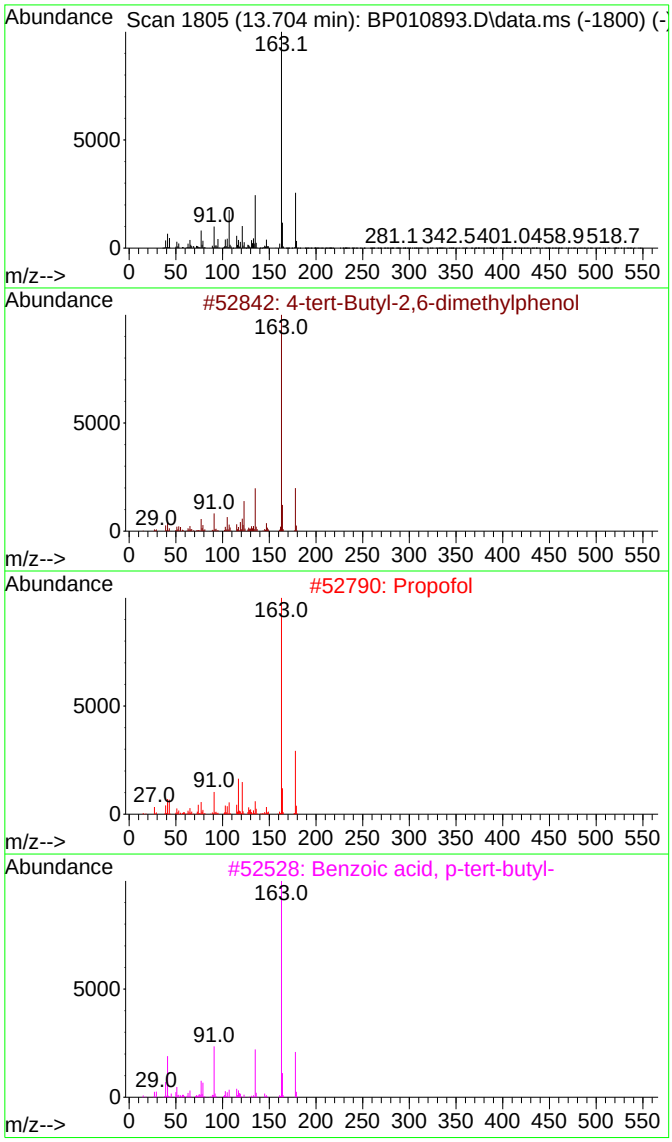
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

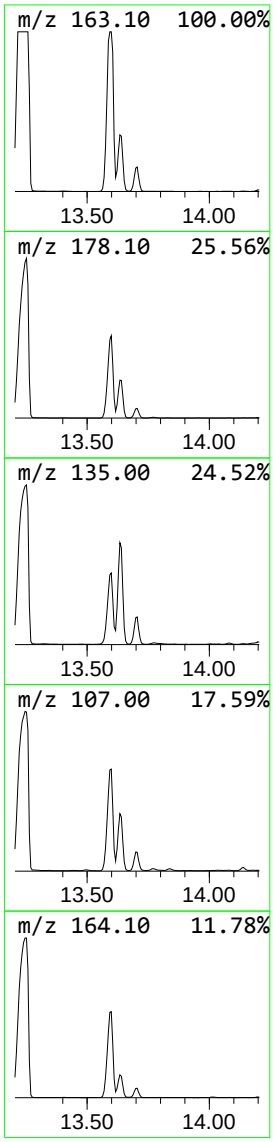
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 4-tert-Butyl-2,6-dimethylph... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.704	25.53 ng/ul	5881370	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-tert-Butyl-2,6-dimethylphenol	178	C12H18O	000879-97-0	90
2	Propofol	178	C12H18O	002078-54-8	81
3	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	81
4	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	74
5	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	72





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Misc :
ALS Vial : 26 Sample Multiplier: 1

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BNA_P
ClientSampleId :
EW4S6RE

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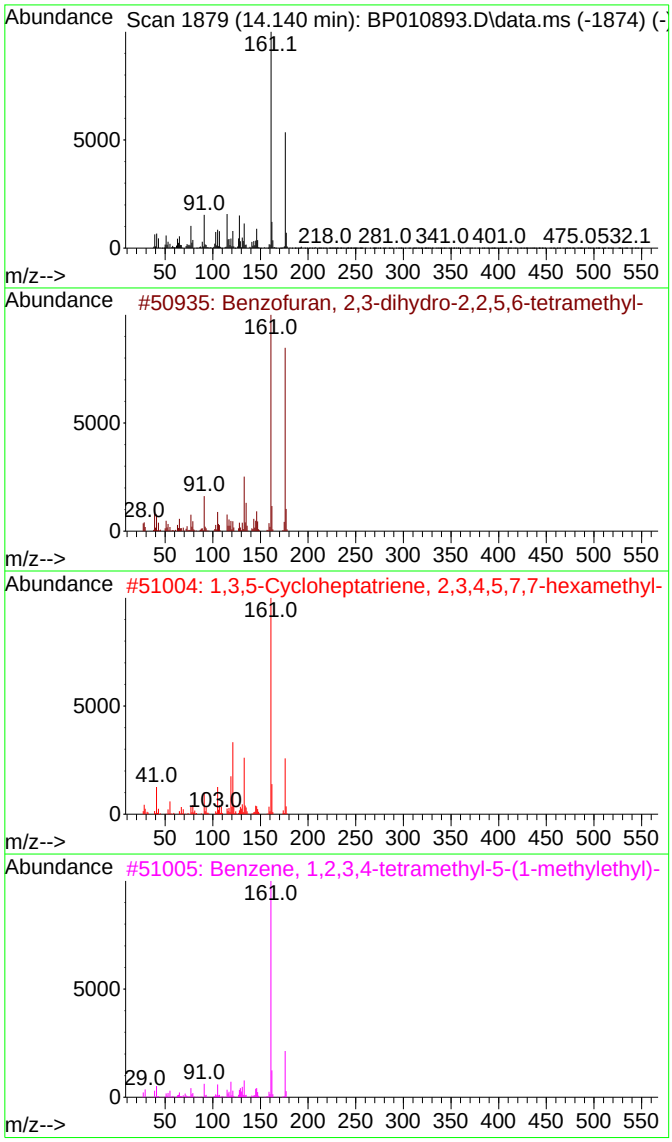
Quant Title : SVOA CALIBRATION

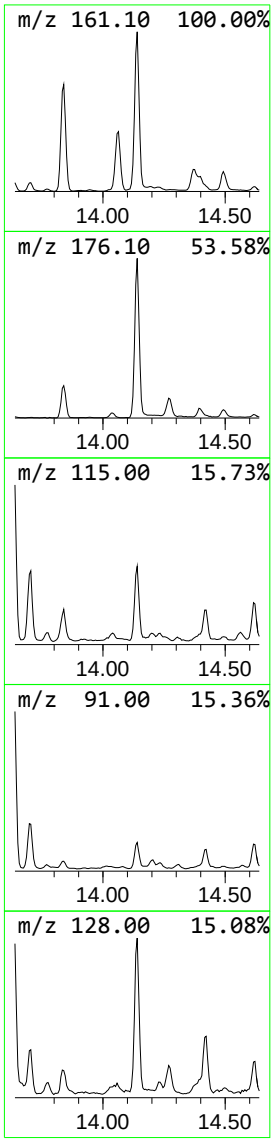
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzofuran, 2,3-dihydro-2,2... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.140	11.81 ng/ul	2719800	Acenaphthene-d10	14.369

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	1,3,5-Cycloheptatriene, 2,3,4,5,...	176	C13H20	074779-68-3	80
3	Benzene, 1,2,3,4-tetramethyl-5-(...	176	C13H20	061142-67-4	72
4	Benzofuran, 2,3-dihydro-2,2,4,6-...	176	C12H16O	003698-49-5	72
5	(1R)-1-(4-tert-butylphenyl)ethan...	177	C12H19N	089538-65-8	72





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Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

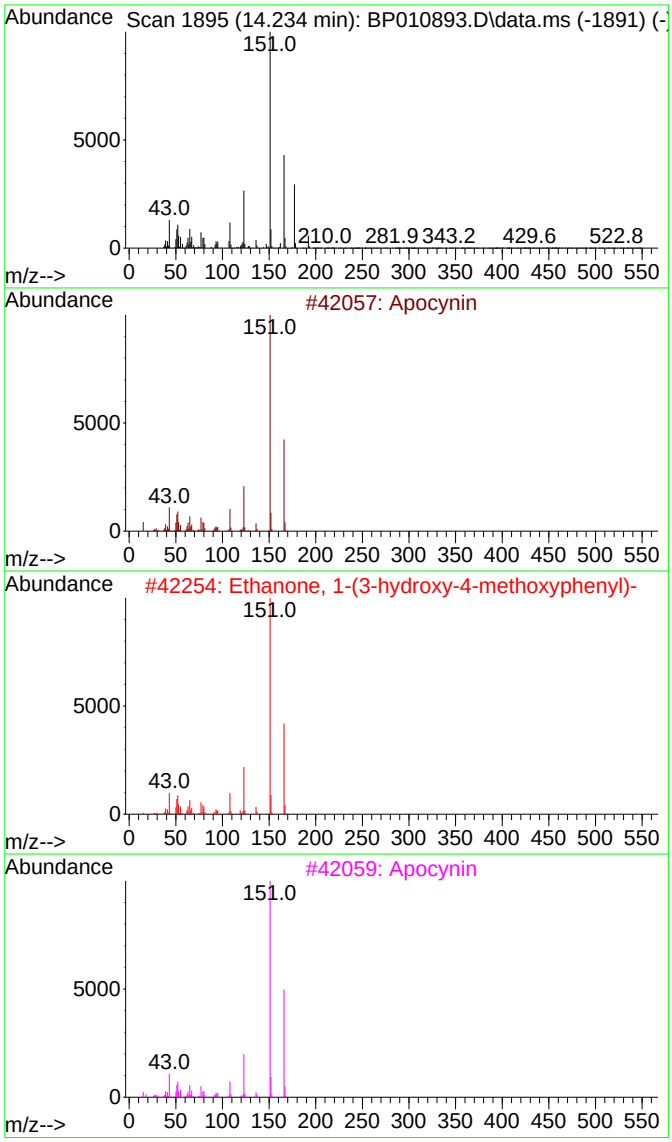
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

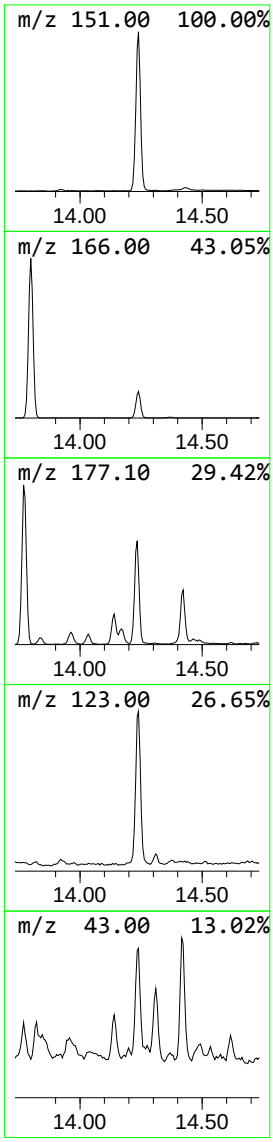
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Apocynin Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.234	8.88 ng/ul	2046210	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Apocynin	166	C9H10O3	000498-02-2	96
2	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	95
3	Apocynin	166	C9H10O3	000498-02-2	95
4	Apocynin	166	C9H10O3	000498-02-2	87
5	Ethanone, 1-(3-hydroxy-4-methoxy...	166	C9H10O3	006100-74-9	87





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

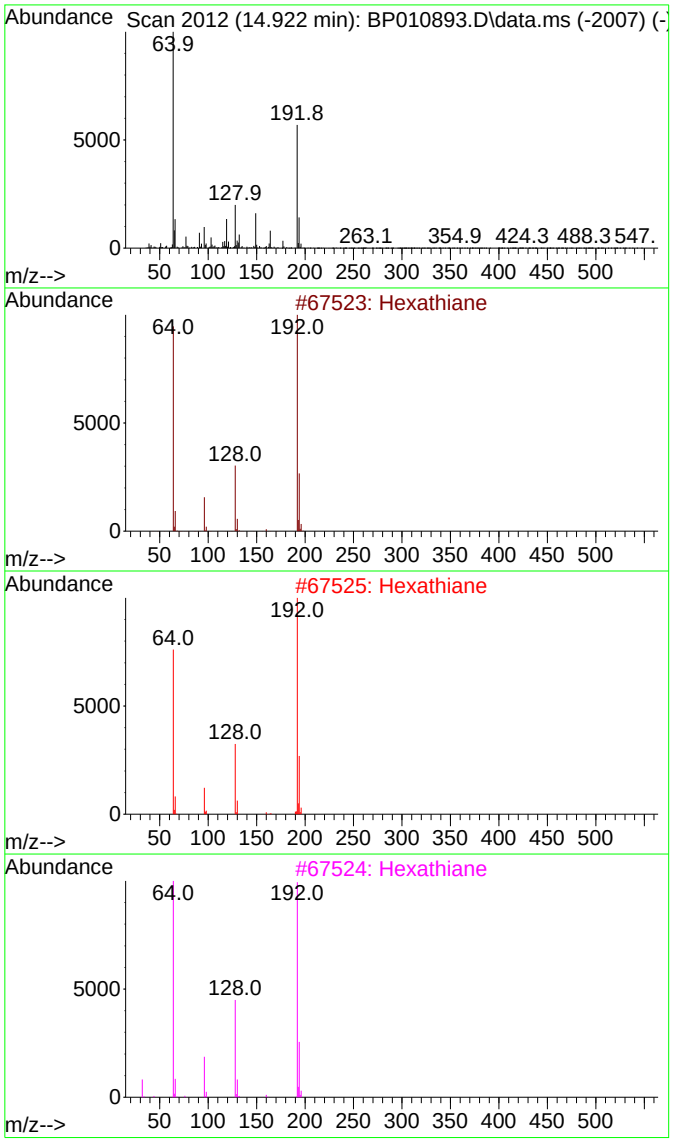
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

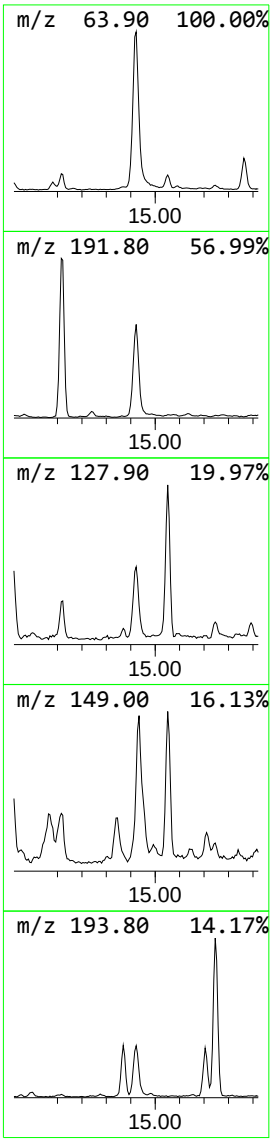
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 Hexathiane Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.922	4.21 ng/ul	968721	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexathiane	192	S6	013798-23-7	91
2	Hexathiane	192	S6	013798-23-7	91
3	Hexathiane	192	S6	013798-23-7	91
4	2-Norbornanone, 1-(epithioethyl)...	228	C11H16O3S	001782-93-0	59
5	Di(tert-butyl) trisulfide, perfl...	534	C8F18S3	016005-64-4	38





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

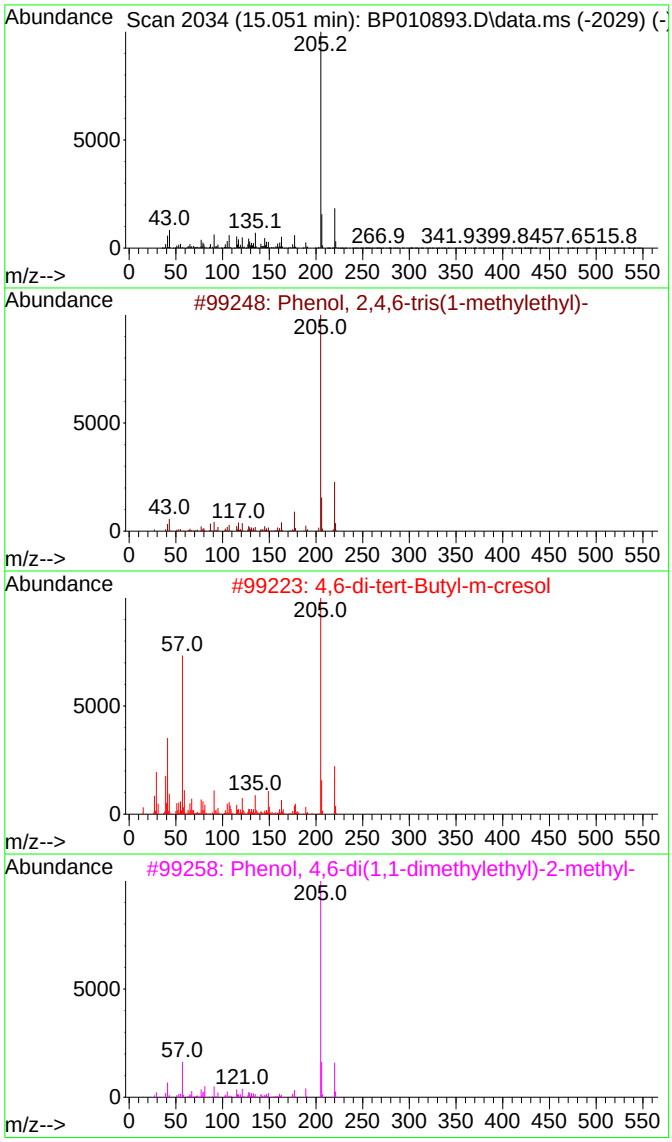
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

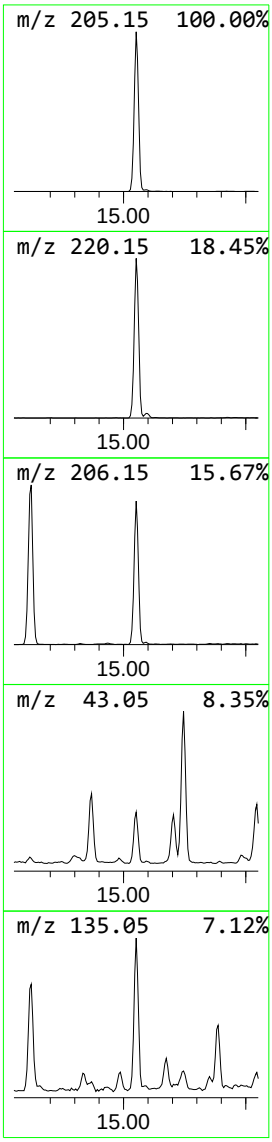
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Phenol, 2,4,6-tris(1-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.051	22.92 ng/ul	5279720	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-tris(1-methylethyl)-	220	C15H24O	002934-07-8	90
2	4,6-di-tert-Butyl-m-cresol	220	C15H24O	000497-39-2	83
3	Phenol, 4,6-di(1,1-dimethylethyl...	220	C15H24O	000616-55-7	81
4	Propanoic acid, 2-methyl-3-[4-t-...	220	C14H20O2	1000131-87-0	81
5	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	81





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

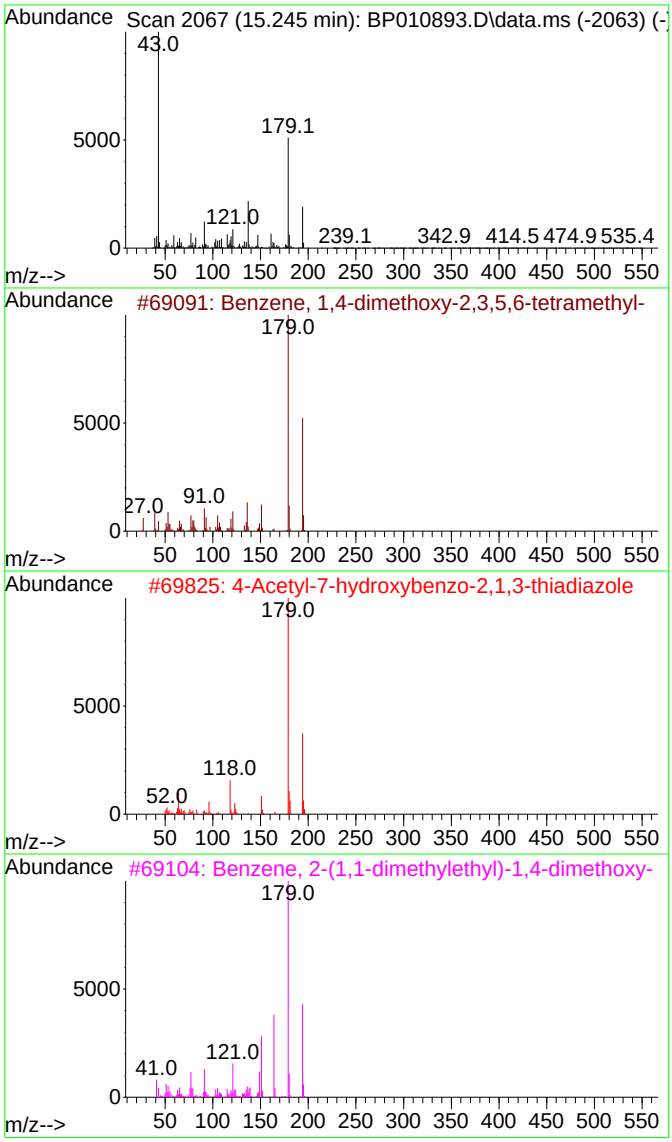
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

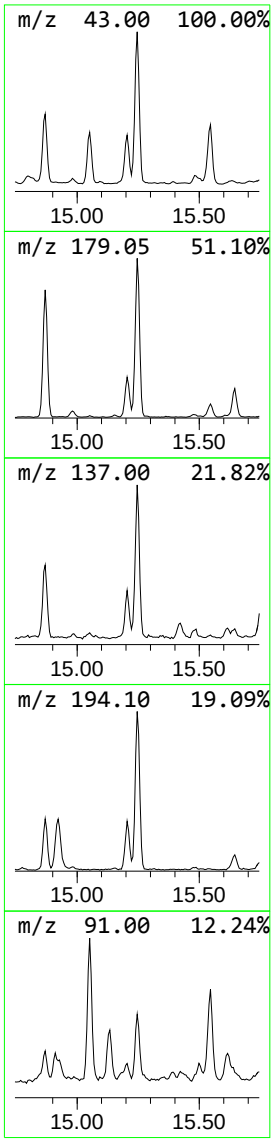
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 Benzene, 1,4-dimethoxy-2,3,... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.246	6.57 ng/ul	1513320	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-dimethoxy-2,3,5,6-t...	194	C12H18O2	013199-54-7	52
2	4-Acetyl-7-hydroxybenzo-2,1,3-th...	194	C8H6N2O2S	120893-38-1	47
3	Benzene, 2-(1,1-dimethylethyl)-1...	194	C12H18O2	021112-37-8	47
4	4-tert-Butylcatechol, dimethyl e...	194	C12H18O2	1000352-79-6	43
5	.alpha.,.alpha.-Diisopropyl-o-me...	222	C14H22O2	010277-90-4	43





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

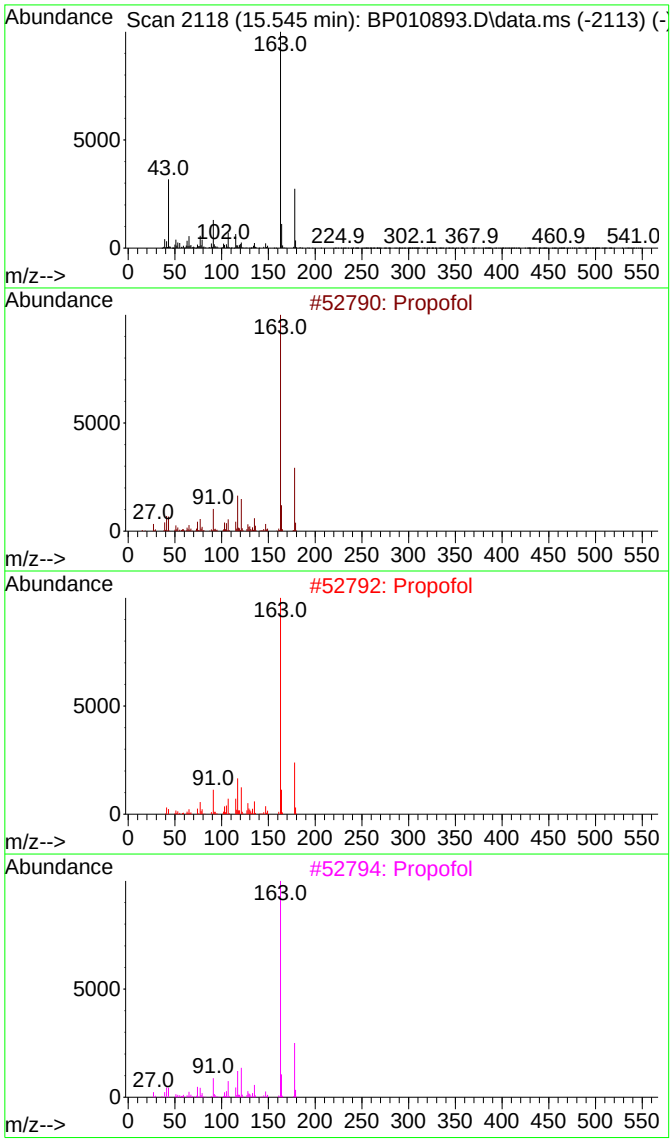
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

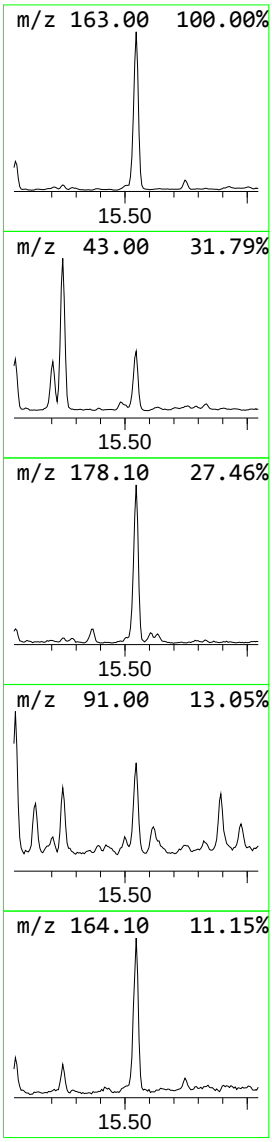
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benzene, 1-(1,1-dimethyleth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.546	7.34 ng/ul	1689840	Acenaphthene-d10	14.369

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	90
2	Propofol	178	C12H18O	002078-54-8	86
3	Propofol	178	C12H18O	002078-54-8	86
4	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	86
5	Propofol	178	C12H18O	002078-54-8	78





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

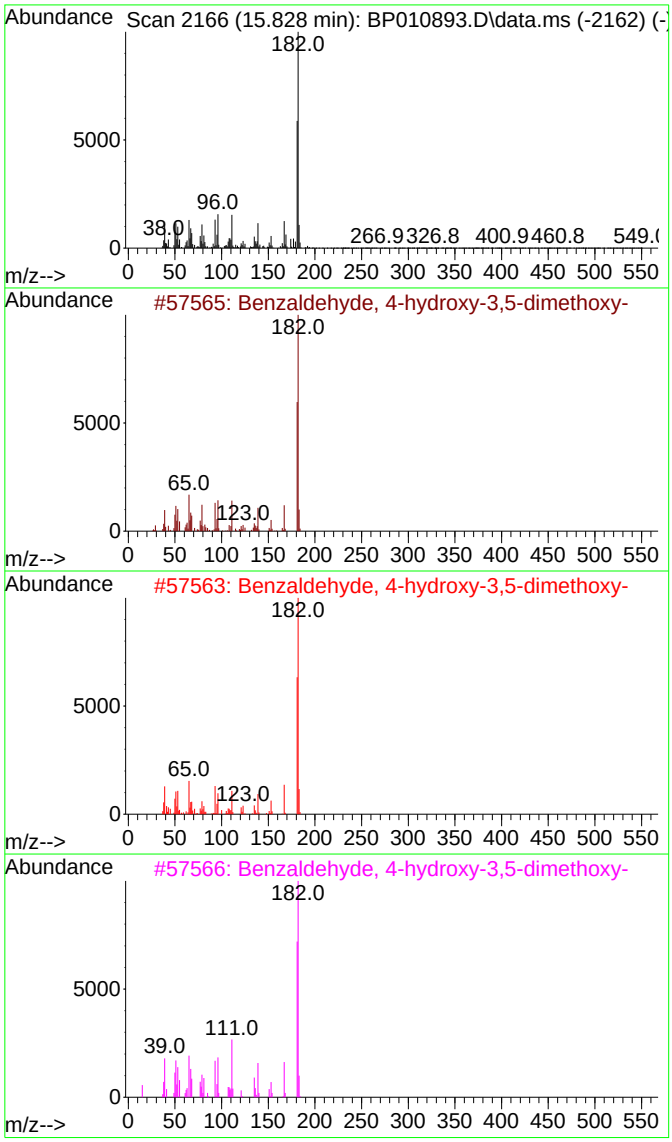
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

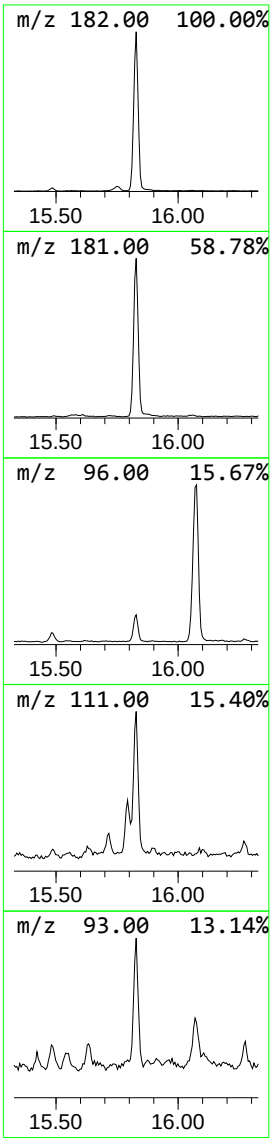
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 Benzaldehyde, 4-hydroxy-3,5... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.828	5.73 ng/ul	1412310	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	98
2	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	98
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	97
5	Benzaldehyde, 4-hydroxy-3,5-dime...	182	C9H10O4	000134-96-3	96





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

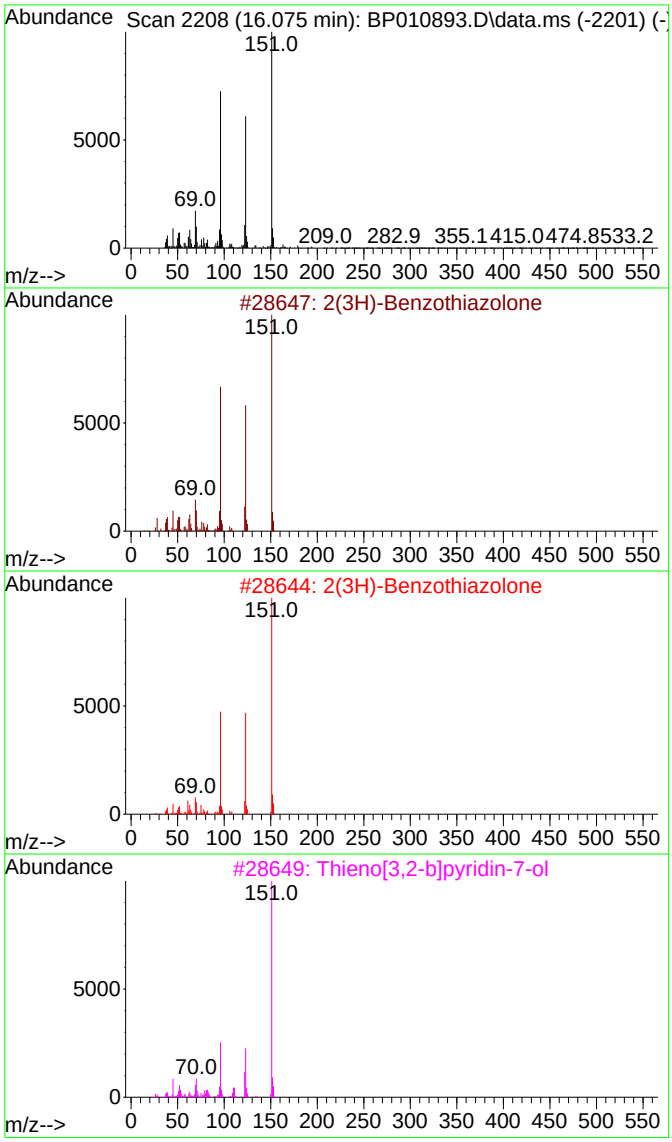
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

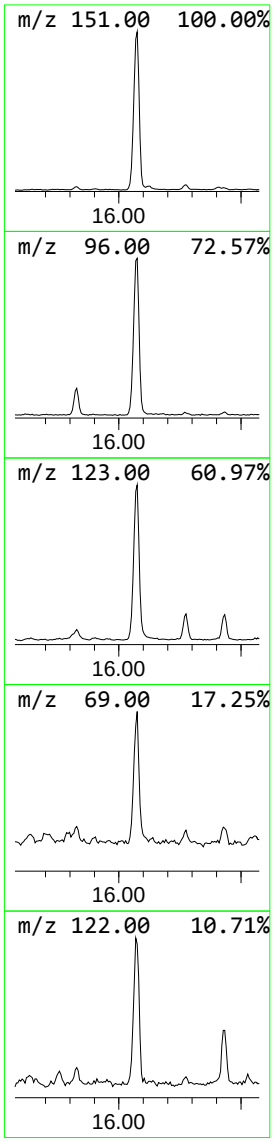
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 2(3H)-Benzothiazolone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	9.59 ng/ul	2362250	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	96
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	58
4	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	53
5	t-Butylhydroquinone	166	C10H14O2	001948-33-0	50





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Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

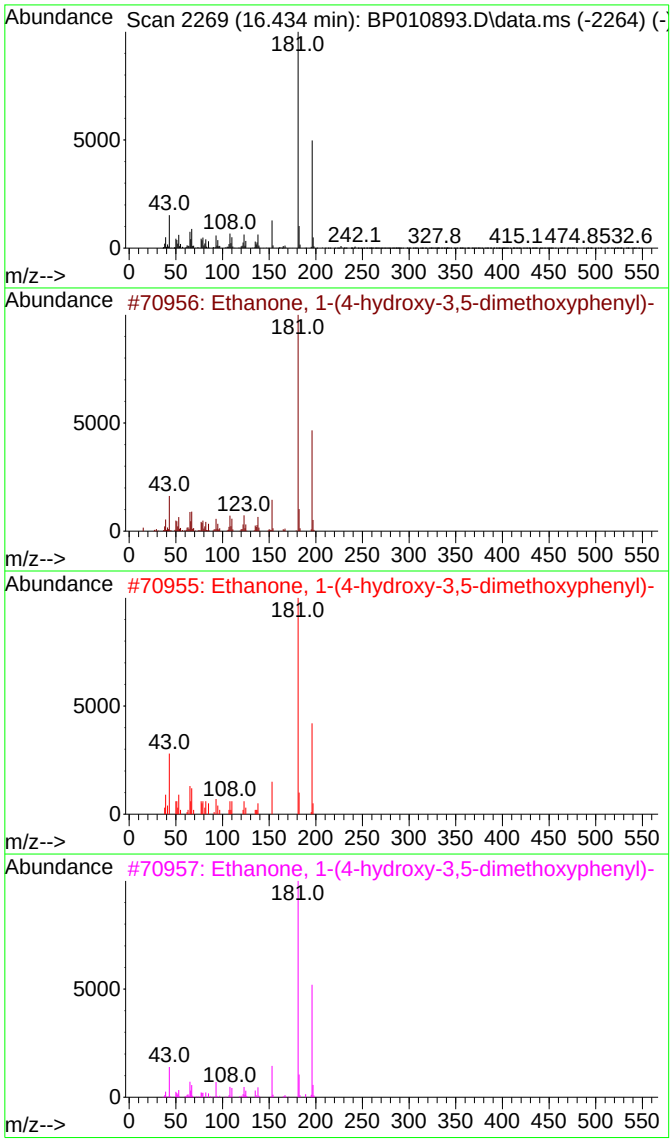
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

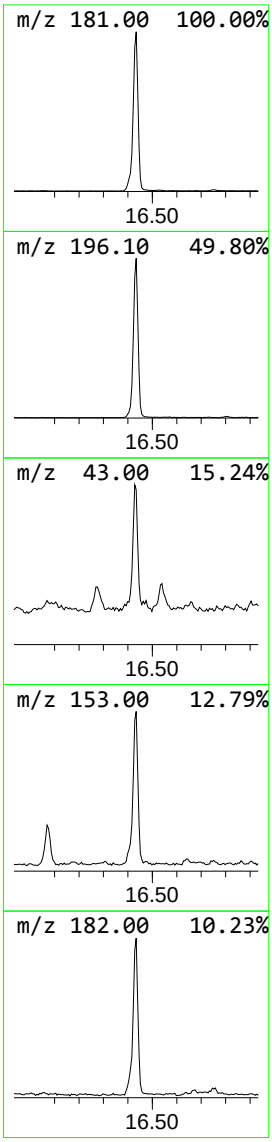
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Ethanone, 1-(4-hydroxy-3,5-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.434	9.61 ng/ul	2368410	Phenanthrene-d10	17.128

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	96
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	90
5	4,5-Dimethoxy-2-hydroxyacetophenone	196	C10H12O4	020628-06-2	87





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Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

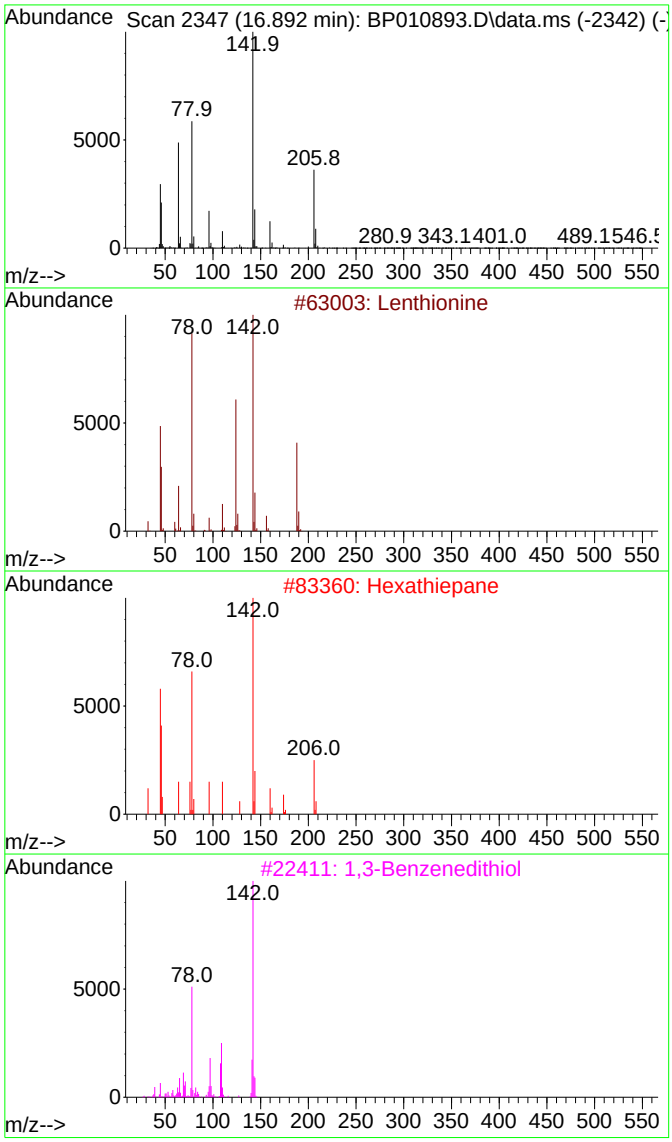
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

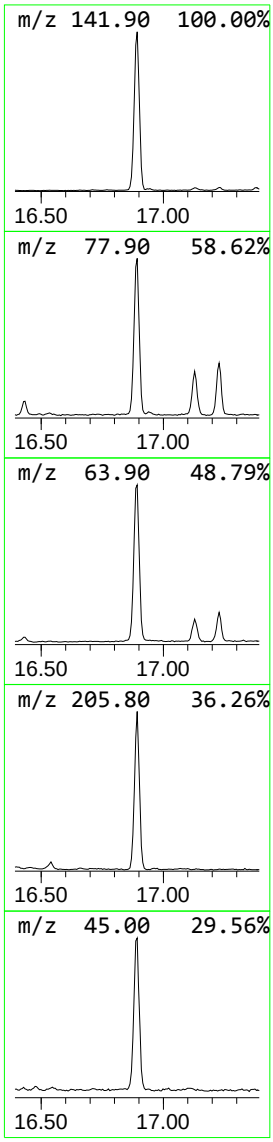
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 unknown-03 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.892	7.28 ng/ul	1793920	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lenthionine	188	C2H4S5	000292-46-6	22
2	Hexathiepane	206	CH2S6	017233-71-5	12
3	1,3-Benzenedithiol	142	C6H6S2	000626-04-0	9
4	4(1H)-Pyrimidinone, 2-(methylthio)-	142	C5H6N2OS	005751-20-2	9
5	Benzenesulfinic-acid-	142	C6H6O2S	000618-41-7	9





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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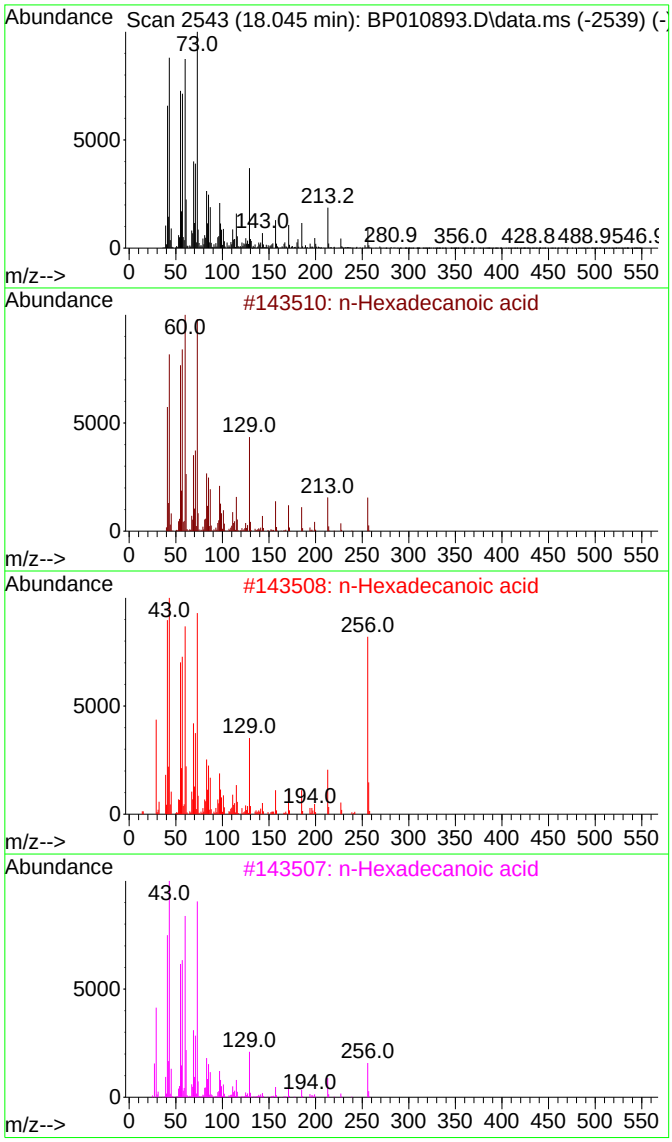
Quant Title : SVOA CALIBRATION

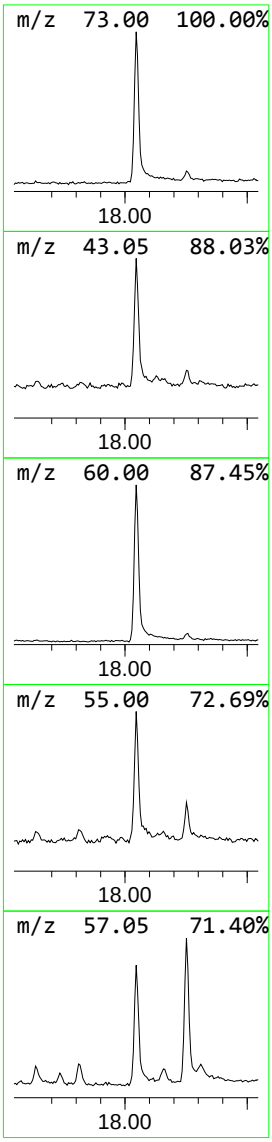
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 31 n-Hexadecanoic acid Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	6.31 ng/ul	1553470	Phenanthrene-d10	17.128

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	95
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Tridecanoic acid	214	C13H26O2	000638-53-9	95





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ClientSampleId :
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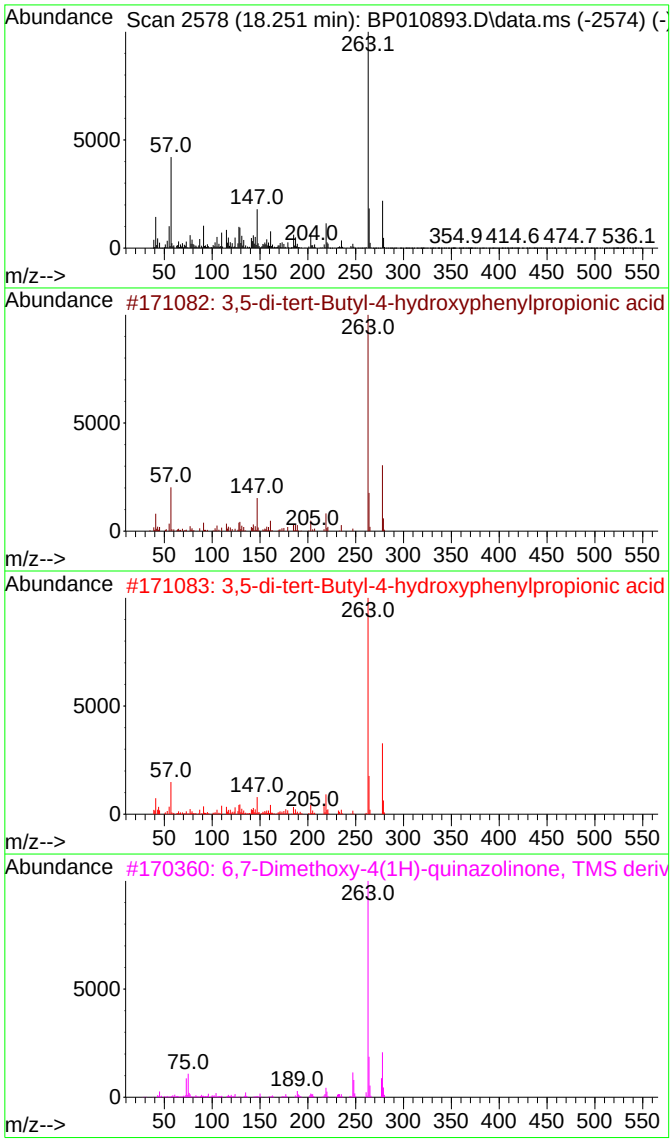
Quant Title : SVOA CALIBRATION

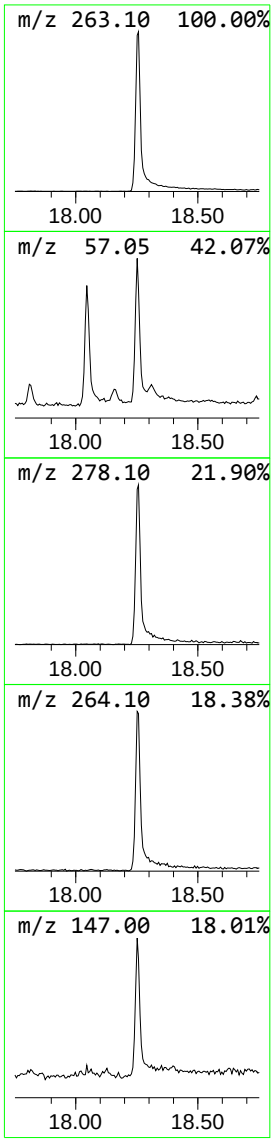
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 32 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	6.20 ng/ul	1526430	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	95
2	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	95
3	6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	64
4	5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	62
5	Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	62





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

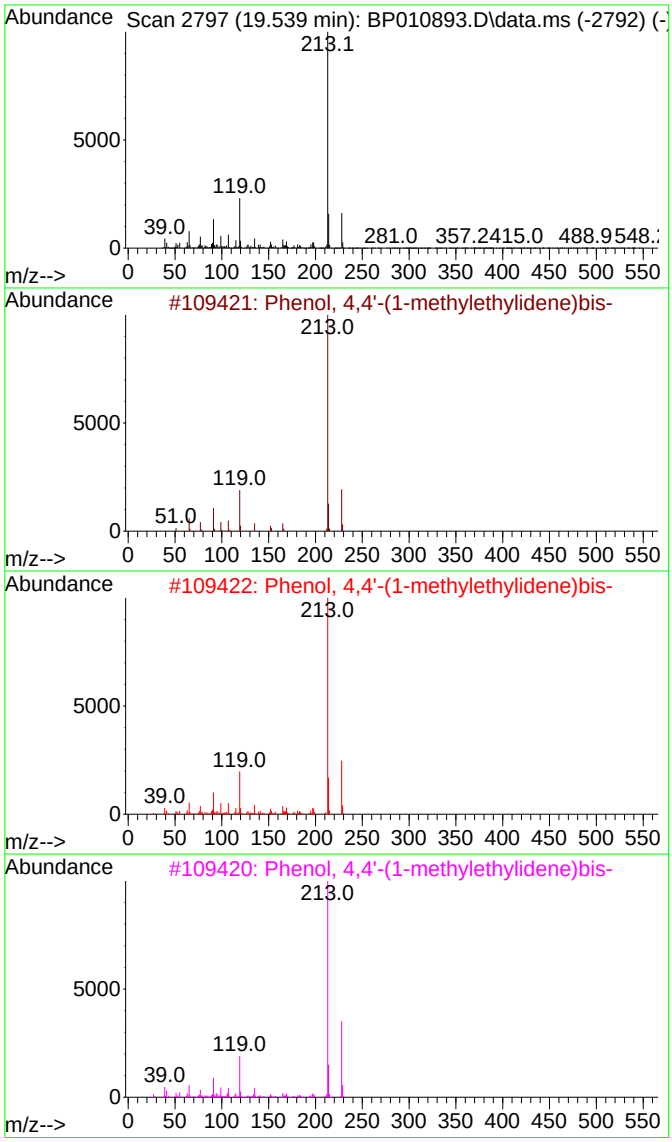
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

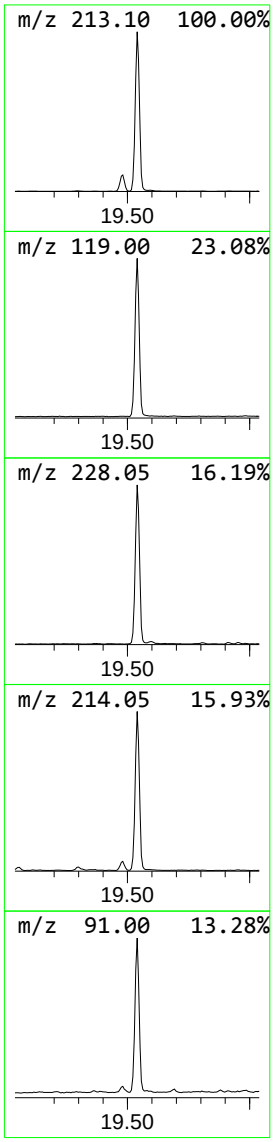
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 34 Phenol, 4,4'-(1-methylethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	46.00 ng/ul	11967200	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	94
4	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	94
5	Diphenolic acid	286	C17H18O4	000126-00-1	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

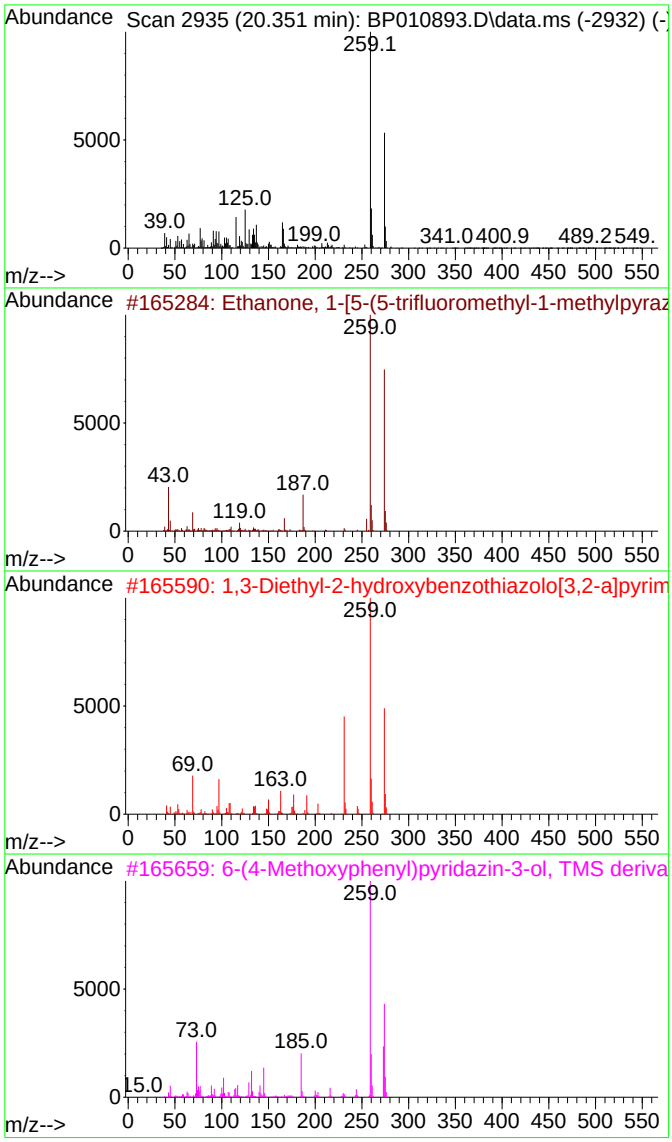
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

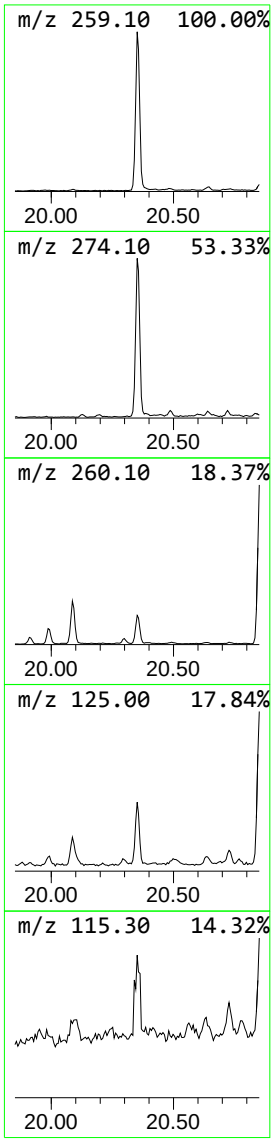
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 36 Ethanone, 1-[5-(5-trifluoro... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.351	4.57 ng/ul	1190130	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-[5-(5-trifluoromethy...	274	C11H9F3N2OS	1000266-44-5	76
2	1,3-Diethyl-2-hydroxybenzothiazo...	274	C14H14N2O2S	1000227-15-3	64
3	6-(4-Methoxyphenyl)pyridazin-3-o...	274	C14H18N2O2Si	1000476-89-3	64
4	Xanthotoxol, trimethylsilyl ether	274	C14H14O4Si	1000462-10-7	64
5	7H-Furo[3,2-g][1]benzopyran-7-on...	274	C15H14O5	054087-32-0	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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

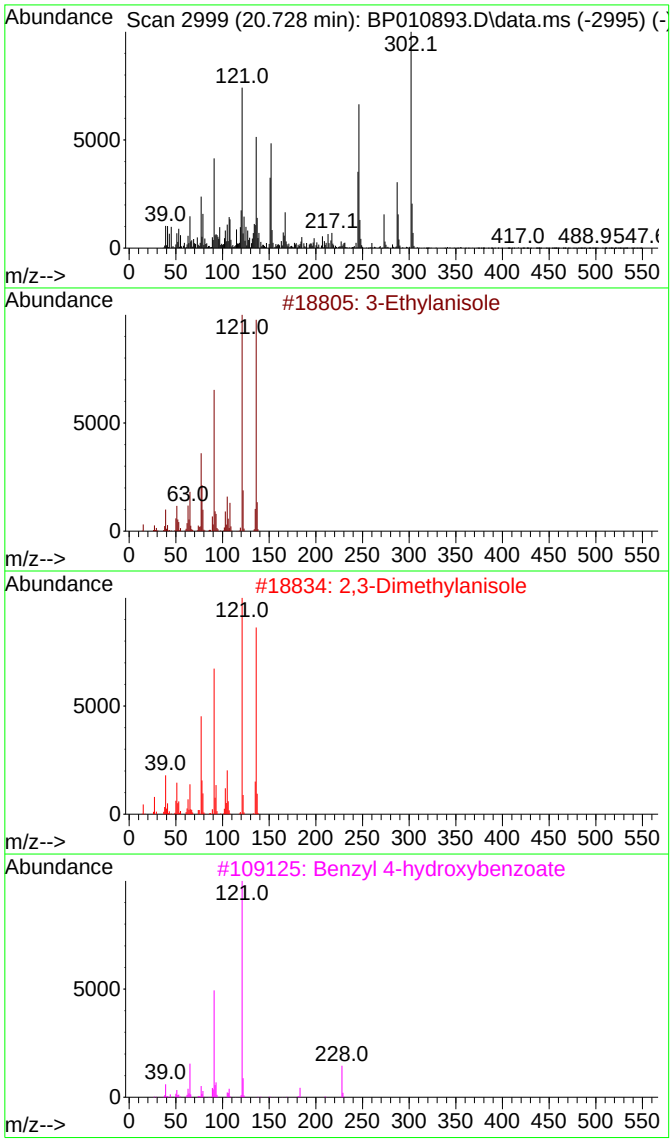
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

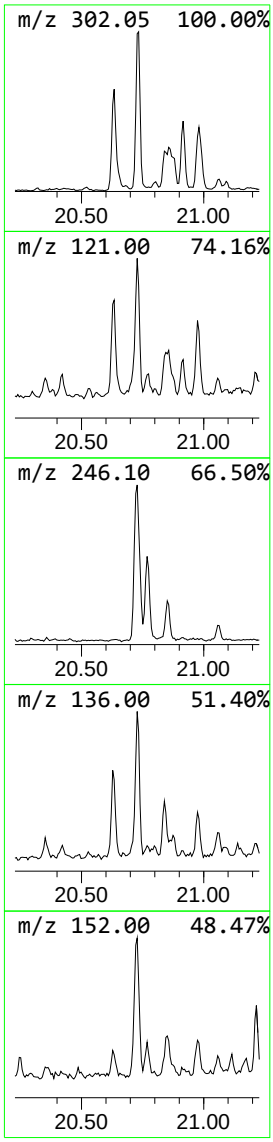
****

Peak Number 38 unknown-04 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.728	4.29 ng/ul	1115690	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Ethylanisole	136	C9H12O	010568-38-4	48
2	2,3-Dimethylanisole	136	C9H12O	002944-49-2	47
3	Benzyl 4-hydroxybenzoate	228	C14H12O3	000094-18-8	25
4	1H-Benzo[b]1,4-diazepin-2(3H)-on...	303	C12H12F3N3O3	147594-83-0	25
5	3,4-Dimethylanisole	136	C9H12O	004685-47-6	20





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

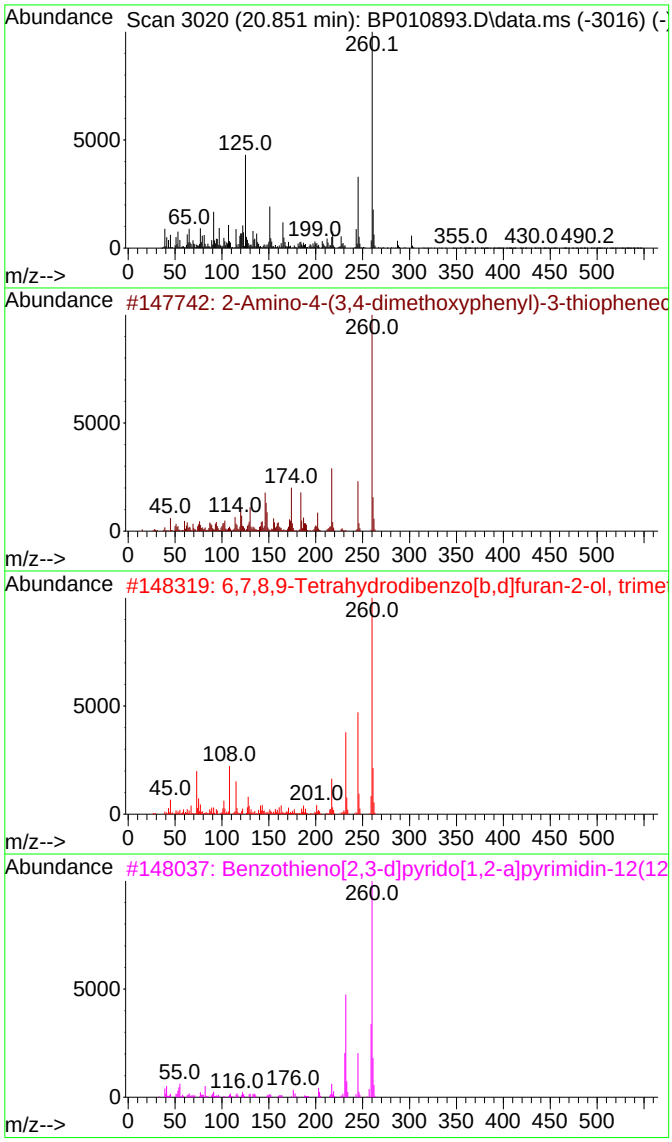
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

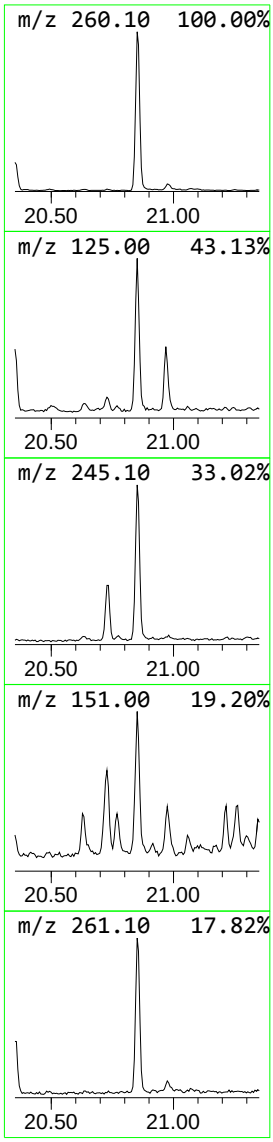
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 unknown-05 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.851	8.27 ng/ul	2150890	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Amino-4-(3,4-dimethoxyphenyl)-...	260	C13H12N2O2S	1000475-04-7	49
2	6,7,8,9-Tetrahydrodibenzo[b,d]fu...	260	C15H20O2Si	1000471-87-6	47
3	Benzothieno[2,3-d]pyrido[1,2-a]p...	260	C14H16N2OS	056929-65-8	46
4	Cyperine	260	C15H16O4	033716-82-4	46
5	1,4-Benzenediamine, N,N'-diphenyl-	260	C18H16N2	000074-31-7	41





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

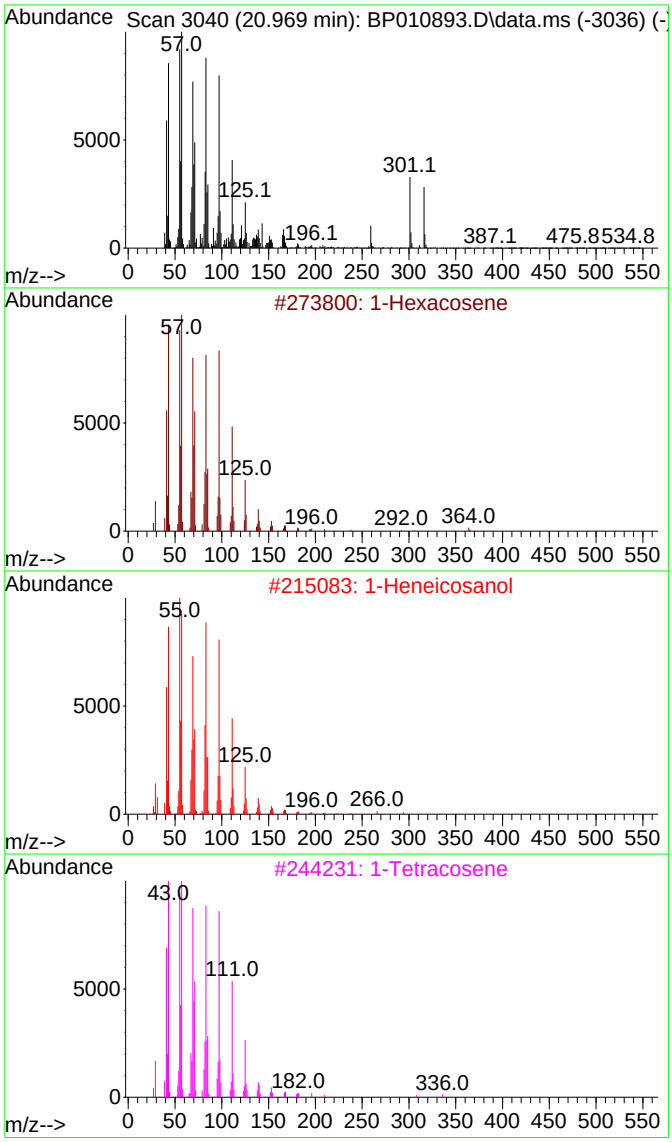
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

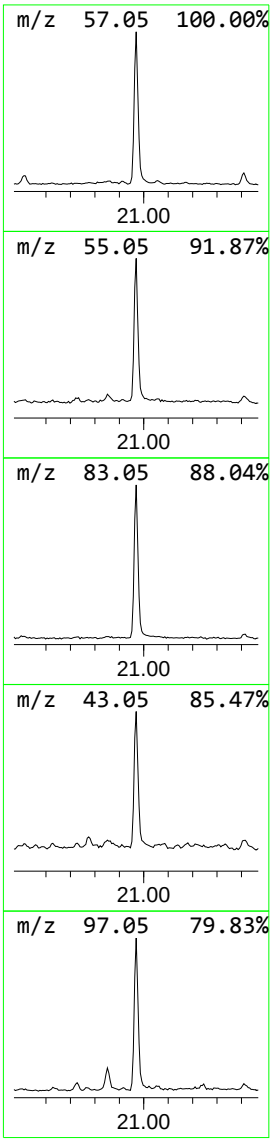
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.969	13.20 ng/ul	3433240	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene	364	C26H52	018835-33-1	93
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	1-Tetracosene	336	C24H48	010192-32-2	93
4	1-Octadecanol	270	C18H38O	000112-92-5	93
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

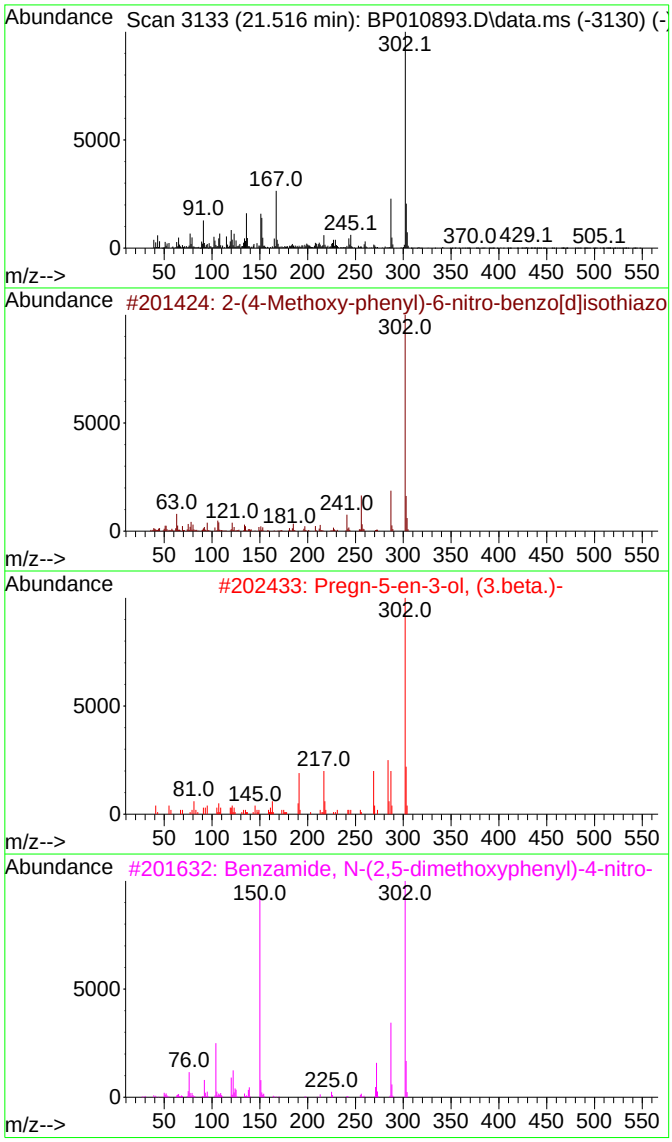
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

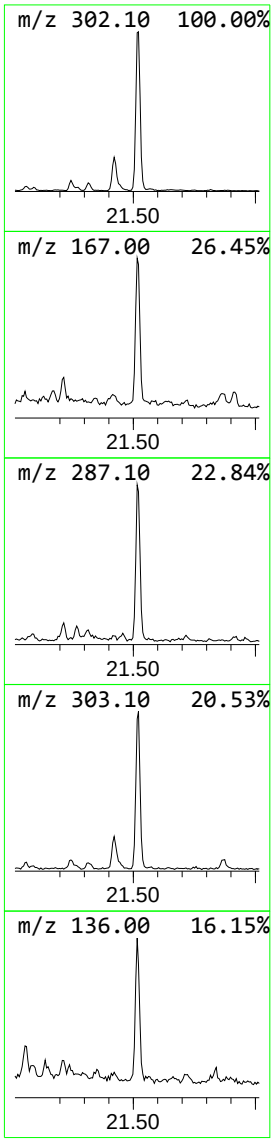
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 2-(4-Methoxy-phenyl)-6-nitr... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.516	4.55 ng/ul	1185040	Chrysene-d12	21.239

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-Methoxy-phenyl)-6-nitro-ben...	302	C14H10N2O4S	1000295-00-9	70
2	Pregn-5-en-3-ol, (3.beta.)-	302	C21H34O	002862-58-0	58
3	Benzamide, N-(2,5-dimethoxypheny...	302	C15H14N2O5	1000307-35-4	50
4	1,1'-Biphenyl, 2,3,4,4'-tetramet...	302	C18H22O4	123705-88-4	50
5	Retinoyl fluoride (all-trans)	302	C20H27FO	083802-77-1	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010893.D
Acq On : 28 Jun 2022 00:27
Operator : CG/JU
Sample : N3374-01RE
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW4S6RE

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	6.228	16.2	ng/ul	2510850	1	7.711	3105540	20.0
1-Hexanol, 2-et...	7.799	25.4	ng/ul	3939340	1	7.711	3105540	20.0
unknown-01	7.958	9.2	ng/ul	1422950	1	7.711	3105540	20.0
3,5-Dimethylhex...	8.558	4.1	ng/ul	637285	1	7.711	3105540	20.0
unknown-02	9.016	5.3	ng/ul	824288	1	7.711	3105540	20.0
p-Cumenol	10.905	5.9	ng/ul	80383700	2	10.522	271411000	20.0
Phenol, p-tert-...	11.963	20.3	ng/ul	275721000	2	10.522	271411000	20.0
(DEL) Alkane: C...	12.516	3.5	ng/ul	816372	3	14.369	4607030	20.0
(DEL) Alkane: C...	12.569	3.0	ng/ul	697827	3	14.369	4607030	20.0
Propofol	12.763	31.8	ng/ul	7317480	3	14.369	4607030	20.0
Phenol, 2,4-bis...	13.252	630.3	ng/ul	145193000	3	14.369	4607030	20.0
Vanillin	13.340	12.3	ng/ul	2831070	3	14.369	4607030	20.0
Phenol, 2,5-bis...	13.599	236.9	ng/ul	54562400	3	14.369	4607030	20.0
Phenol, 3,5-bis...	13.634	84.3	ng/ul	19428100	3	14.369	4607030	20.0
4-tert-Butyl-2,...	13.704	25.5	ng/ul	5881370	3	14.369	4607030	20.0
Benzofuran, 2,3...	14.140	11.8	ng/ul	2719800	3	14.369	4607030	20.0
Apocynin	14.234	8.9	ng/ul	2046210	3	14.369	4607030	20.0
Hexathiane	14.922	4.2	ng/ul	968721	3	14.369	4607030	20.0
Phenol, 2,4,6-t...	15.051	22.9	ng/ul	5279720	3	14.369	4607030	20.0
Benzene, 1,4-di...	15.246	6.6	ng/ul	1513320	3	14.369	4607030	20.0
Benzene, 1-(1,1...	15.546	7.3	ng/ul	1689840	3	14.369	4607030	20.0
Benzaldehyde, 4...	15.828	5.7	ng/ul	1412310	4	17.128	4926640	20.0
2(3H)-Benzothia...	16.075	9.6	ng/ul	2362250	4	17.128	4926640	20.0
Ethanone, 1-(4-...	16.434	9.6	ng/ul	2368410	4	17.128	4926640	20.0
unknown-03	16.892	7.3	ng/ul	1793920	4	17.128	4926640	20.0
n-Hexadecanoic ...	18.045	6.3	ng/ul	1553470	4	17.128	4926640	20.0
3,5-di-tert-But...	18.251	6.2	ng/ul	1526430	4	17.128	4926640	20.0
Phenol, 4,4'-(1...	19.539	46.0	ng/ul	11967200	5	21.239	5203280	20.0
Ethanone, 1-[5-...	20.351	4.6	ng/ul	1190130	5	21.239	5203280	20.0
unknown-04	20.728	4.3	ng/ul	1115690	5	21.239	5203280	20.0
unknown-05	20.851	8.3	ng/ul	2150890	5	21.239	5203280	20.0
(DEL) Alkane: S...	20.969	13.2	ng/ul	3433240	5	21.239	5203280	20.0
2-(4-Methoxy-ph...	21.516	4.5	ng/ul	1185040	5	21.239	5203280	20.0