LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

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-BP111023.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP018237.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	10	13	18	rVB2	13818	17972	1.23%	0.138%
2	3.605	85	88	98	rBV	59811	114526	7.84%	0.882%
3	3.769	114	116	122	rVB2	5203	6088	0.42%	0.047%
4	4.452	229	232	235	rVB5	1985	2333	0.16%	0.018%
5	5.016	326	328	334	rBV6	1111	1809	0.12%	0.014%
6	5.463	402	404	406	rVV3	2250	2017	0.14%	0.016%
7	6.134	513	518	521	rBV6	1782	2962	0.20%	0.023%
8	6.340	550	553	555	rVB3	1437	1754	0.12%	0.014%
9	7.040	668	672	674	rBV5	2167	2893	0.20%	0.022%
10	7.134	683	688	703	rBV	123966	230023	15.75%	1.772%
11	7.340	719	723	737	rVB2	8370	21896	1.50%	0.169%
12	7.675	779	780	786	rVB6	1982	2519	0.17%	0.019%
13	7.781	792	798	810	rBV	132091	264787	18.13%	2.039%
14	8.399	900	903	908	rBV7	1369	1894	0.13%	0.015%
15	8.563	925	931	940	rBV9	7889	24993	1.71%	0.193%
16	8.646	943	945	948	rVB3	2037	2216	0.15%	0.017%
17	8.993	996	1004	1022	rBV	145972	338234	23.16%	2.605%
18	9.140	1026	1029	1030	rBV3	2120	2129	0.15%	0.016%
19	9.746	1131	1132	1136	rBV4	1485	2015	0.14%	0.016%
20	10.293	1220	1225	1231	rBV6	7126	18137	1.24%	0.140%
21	10.557	1266	1270	1271	rBV4	1571	2078	0.14%	0.016%
22	10.598	1271	1277	1291	rVV	179232	332924	22.80%	2.564%
23	10.687	1291	1292	1296	rVB4	2349	2427	0.17%	0.019%
24	10.787	1302	1309	1329	rBV	141176	391559	26.81%	3.016%
25	11.645	1451	1455	1459	rVB7	1679	2482	0.17%	0.019%
26	12.116	1533	1535	1540	rBV4	1421	1819	0.12%	0.014%
27	13.022	1685	1689	1690	rBV4	1632	1915	0.13%	0.015%
28	13.204	1718	1720	1725	rVB6	2227	2460	0.17%	0.019%
29	13.298	1729	1736	1750	rBV	534160	787451	53.92%	6.065%
30	13.575	1781	1783	1788	rBV5	1030	1701	0.12%	0.013%
31	13.669	1795	1799	1802	rBV6	1890	3449	0.24%	0.027%
32	13.745	1808	1812	1814	rBV5	3373	5425	0.37%	0.042%
33	13.881	1830	1835	1853	rVV	413589	654096	44.79%	5.038%
34	13.998	1853	1855	1862	rVB8	1663	2609	0.18%	0.020%
35	14.075	1862	1868	1870	rVB7	1674	2429	0.17%	0.019%
36	14.157	1876	1882	1892	rBV	472591	710419	48.64%	5.472%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

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-BP111023.MA.M
Title : SVOA CALIBRATION

37	14.292	1901	1905	1910	rVB7	2146	3159	0.22%	0.024%
38	14.386	1920	1921	1926	rVB4	1498	2141	0.15%	0.016%
39	14.463	1926	1934	1946	rBV	281711	445623	30.51%	3.432%
40	15.110	2041	2044	2047	rVV4	2708	3066	0.21%	0.024%
41	15.192	2055	2058	2064	rBV8	1908	2824	0.19%	0.022%
42	15.251	2064	2068	2072	rVB6	2944	4683	0.32%	0.036%
43	15.333	2079	2082	2083	rVV3	2950	2825	0.19%	0.022%
44	15.345	2083	2084	2089	rVB5	2555	2439	0.17%	0.019%
45	15.392	2089	2092	2095	rBV5	1287	1807	0.12%	0.014%
46	15.463	2098	2104	2125	rBV	644640	1010889	69.22%	7.786%
47	15.616	2128	2130	2133	rVV4	2485	2514	0.17%	0.019%
48	15.669	2136	2139	2142	rBV4	1822	2407	0.16%	0.019%
49	15.722	2142	2148	2149	rBV6	2587	3844	0.26%	0.030%
50	15.739	2149	2151	2154	rVB4	2642	2380	0.16%	0.018%
51	15.881	2166	2175	2179	rVV	12684	21086	1.44%	0.162%
52	15.922	2179	2182	2186	rVV2	9963	14013	0.96%	0.108%
53	15.963	2186	2189	2194	rVV6	3518	6427	0.44%	0.050%
54	16.180	2220	2226	2232	rBV	30079	40833	2.80%	0.315%
55	16.275	2236	2242	2246	rVV	97209	129053	8.84%	0.994%
56	16.333	2246	2252	2258	rVV2	120799	248530	17.02%	1.914%
57	16.398	2258	2263	2267	rVV2	98438	175998	12.05%	1.356%
58	16.439	2267	2270	2275	rVV	59377	81787	5.60%	0.630%
59	16.498	2275	2280	2281	rVV2	30817	46060	3.15%	0.355%
60	16.516	2281	2283	2285	rVV	29027	34295	2.35%	0.264%
61	16.545	2285	2288	2291	rVV	32245	40718	2.79%	0.314%
62	16.598	2291	2297	2303	rVV4	70818	139261	9.54%	1.073%
63	16.669	2303	2309	2314	rVV	97671	150108	10.28%	1.156%
64	16.710	2314	2316	2318	rVV3	27075	32208	2.21%	0.248%
65	16.739	2318	2321	2330	rVB2	41334	63989	4.38%	0.493%
66	16.810	2330	2333	2334	rBV3	1617	1681	0.12%	0.013%
67	16.886	2339	2346	2348	rBV7	3062	5211	0.36%	0.040%
68	16.998	2361	2365	2368	rBV5	3250	4725	0.32%	0.036%
69	17.086	2376	2380	2382	rVB5	2542	2680	0.18%	0.021%
70	17.245	2401	2407	2415	rBV	374547	541039	37.05%	4.167%
71	17.345	2418	2424	2441	rVV	730783	1147122	78.55%	8.835%
72	17.469	2443	2445	2447	rVV2	2679	3444	0.24%	0.027%
73	17.492	2447	2449	2456	rVV8	3945	7388	0.51%	0.057%
74	17.575	2460	2463	2465	rBV4	1958	2128	0.15%	0.016%
75	17.663	2476	2478	2482	rVB5	1755	2070	0.14%	0.016%
76	17.704	2482	2485	2488	rVB4	1879	1780	0.12%	0.014%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

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-BP111023.MA.M
Title : SVOA CALIBRATION

77	17.739	2488	2491	2492	rBV3	1770	2357	0.16%	0.018%
78	17.869	2506	2513	2519	rBV3	9648	16329	1.12%	0.126%
79	18.004	2534	2536	2541	rBV6	1399	2319	0.16%	0.018%
80	18.110	2549	2554	2559	rBV8	4728	9709	0.66%	0.075%
81	18.186	2566	2567	2571	rVB4	2464	2900	0.20%	0.022%
82	18.363	2593	2597	2601	rBV7	3380	5562	0.38%	0.043%
83	18.498	2618	2620	2623	rVB4	2458	2291	0.16%	0.018%
84	18.522	2623	2624	2626	rBV2	2148	1735	0.12%	0.013%
85	18.986	2700	2703	2704	rBV2	4548	4552	0.31%	0.035%
86	19.016	2704	2708	2714	rVB	41281	47193	3.23%	0.363%
87	19.110	2721	2724	2725	rBV3	2357	2194	0.15%	0.017%
88	19.145	2725	2730	2732	rBV6	7893	12076	0.83%	0.093%
89	19.174	2732	2735	2741	rVB3	10882	14080	0.96%	0.108%
90	19.233	2742	2745	2746	rBV2	10389	9371	0.64%	0.072%
91	19.304	2755	2757	2758	rBV2	2557	1692	0.12%	0.013%
92	19.533	2793	2796	2799	rBV5	5107	7483	0.51%	0.058%
93	19.598	2802	2807	2822	rVV	1086740	1460447	100.00%	11.249%
94	20.063	2883	2886	2889	rBV4	10477	11092	0.76%	0.085%
95	20.557	2967	2970	2972	rBV3	8793	8633	0.59%	0.066%
96	21.033	3046	3051	3065	rBV4	70114	185388	12.69%	1.428%
97	21.374	3105	3109	3120	rBV	410168	627320	42.95%	4.832%
98	22.580	3310	3314	3318	rVB	40079	56400	3.86%	0.434%
99	23.680	3494	3501	3511	rBV	680409	1401469	95.96%	10.795%
100	23.851	3523	3530	3545	rVB	292696	689841	47.23%	5.313%

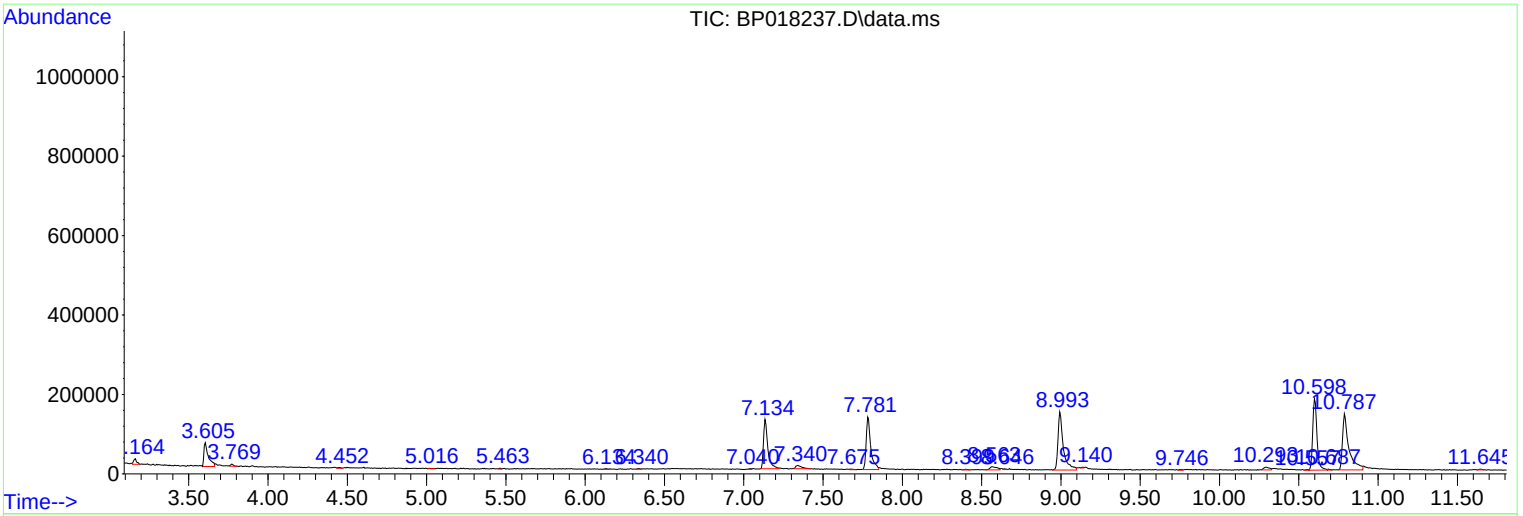
Sum of corrected areas: 12983108

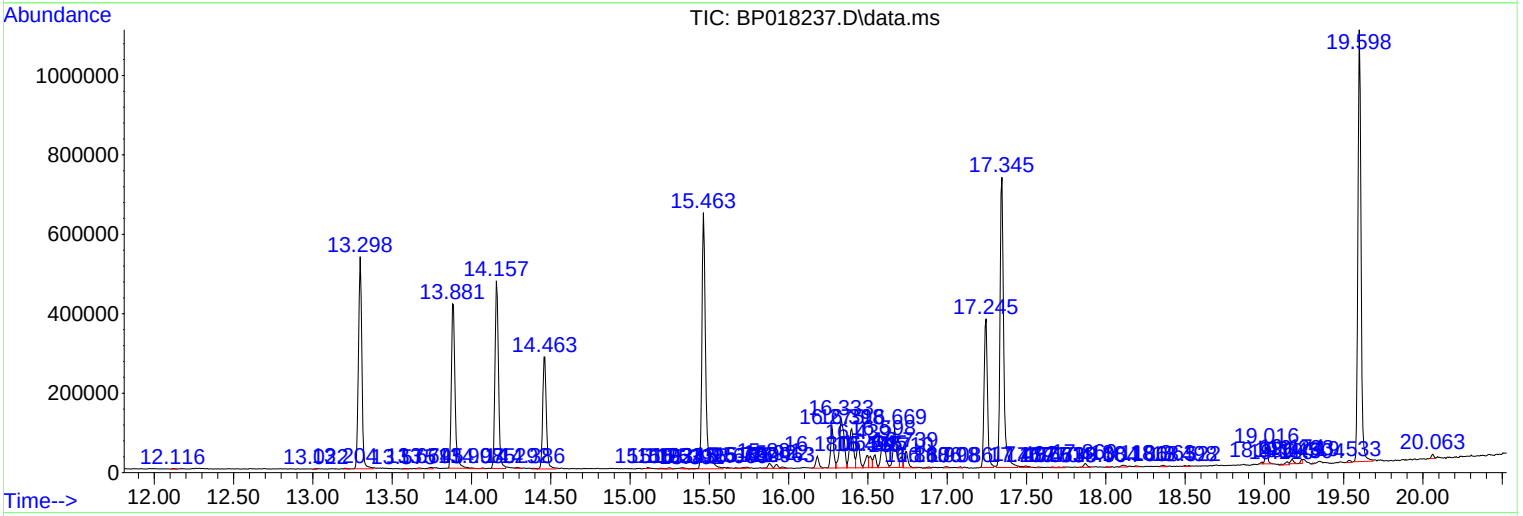
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

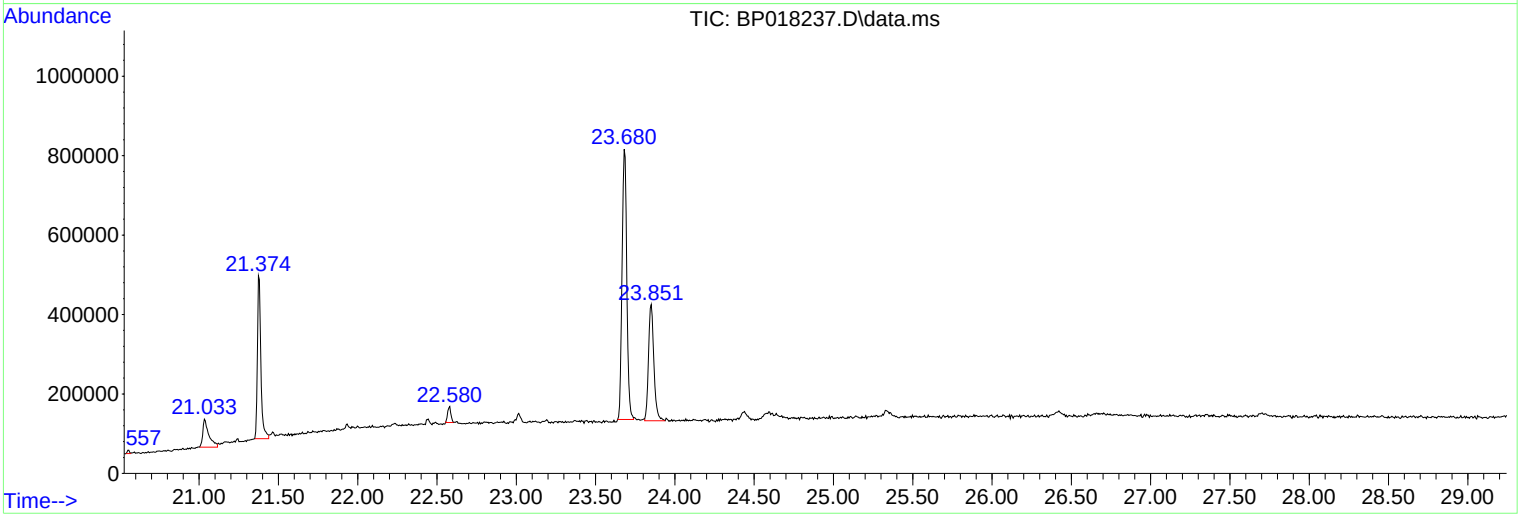
Instrument :
BNA_P
ClientSampleId :
GCDP6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

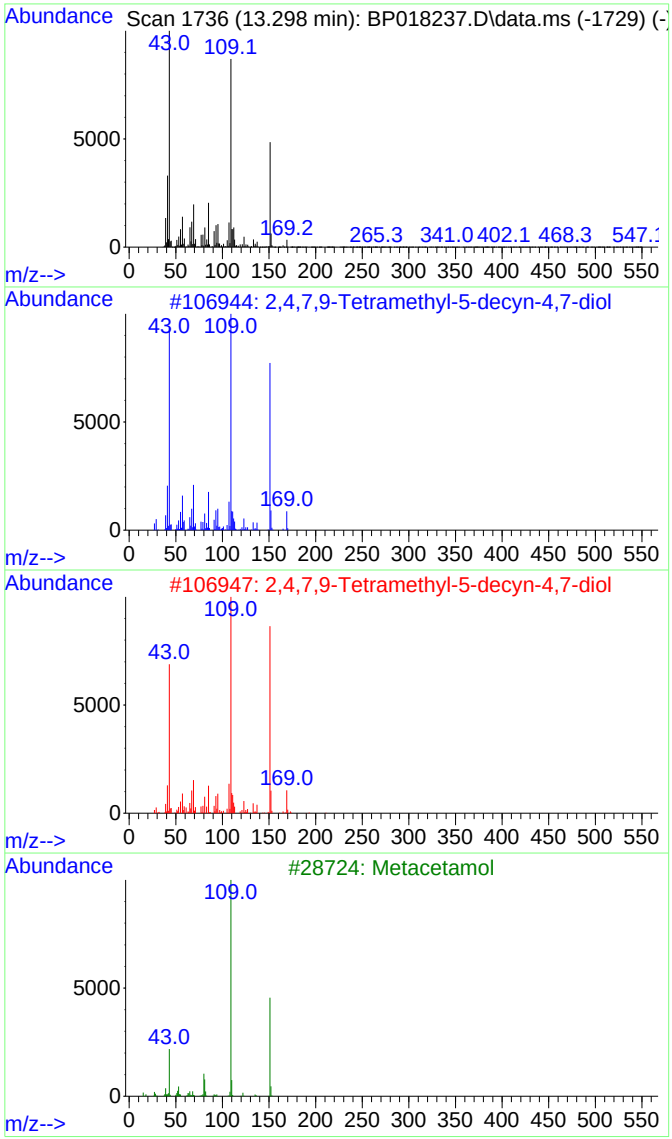
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

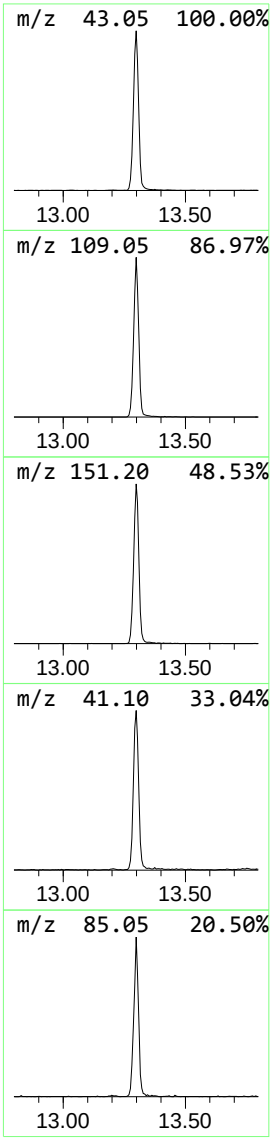
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	35.34 ng/ul	787451	Acenaphthene-d10	14.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	87
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	72
3	Metacetamol	151	C8H9NO2	000621-42-1	52
4	4-Pyridinemethanol, acetate (ester)	151	C8H9NO2	001007-48-3	50
5	Acetaminophen	151	C8H9NO2	000103-90-2	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

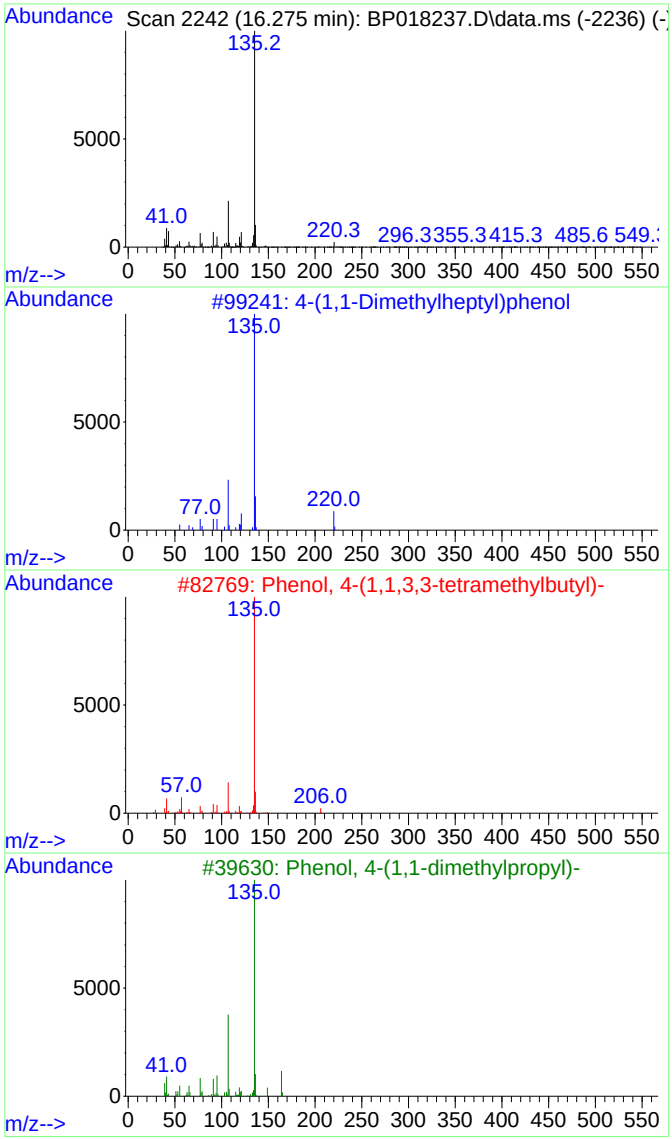
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

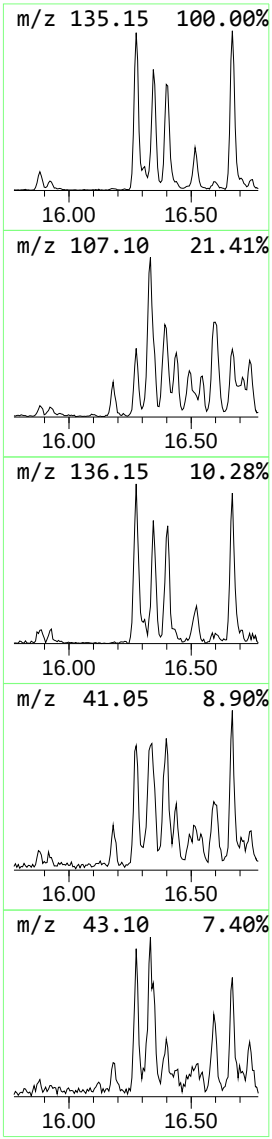
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 4-(1,1-Dimethylheptyl)phenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	4.77 ng/ul	129053	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	87
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	78
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

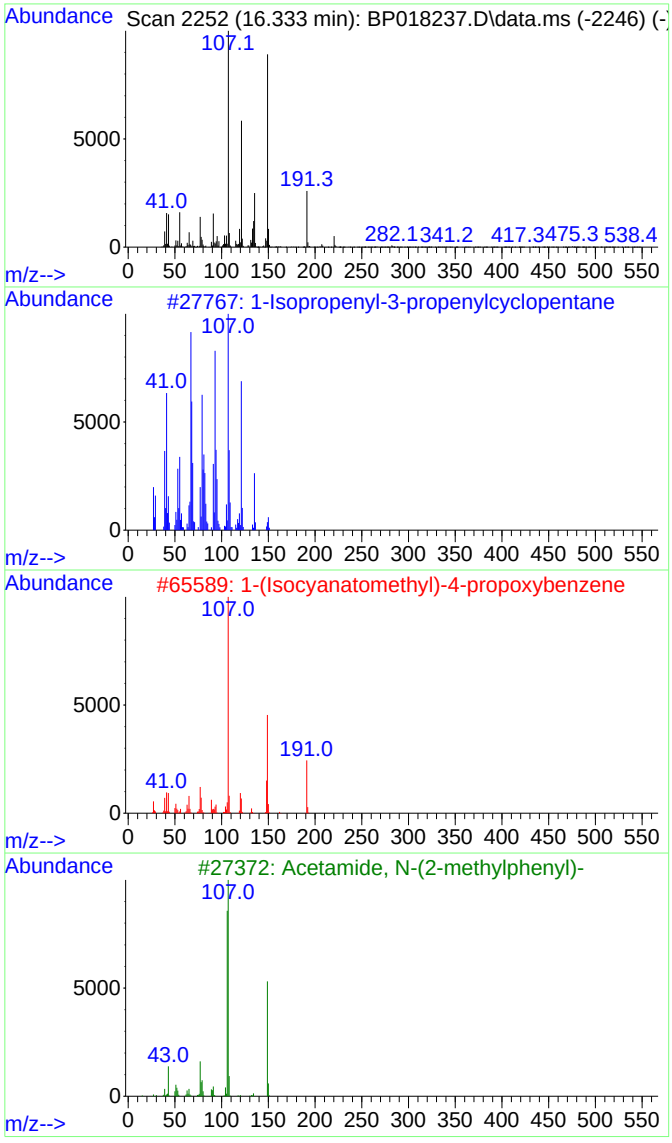
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

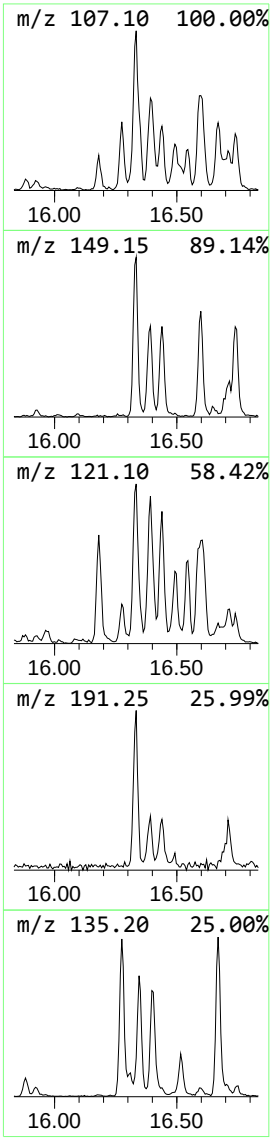
****

Peak Number 3 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.333	9.19 ng/ul	248530	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Isopropenyl-3-propenylcyclopen...	150	C11H18	1000192-81-2	46
2	1-(Isocyanatomethyl)-4-propoxybe...	191	C11H13NO2	936095-07-7	43
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
4	1-(2,6-Dimethyl-4-propoxyphenyl)...	220	C14H20O2	1000190-22-3	38
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

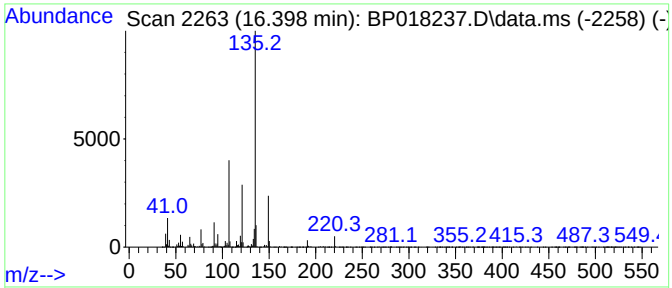
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

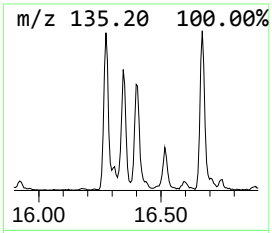
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

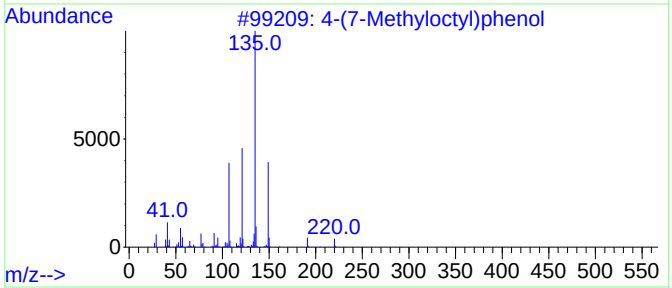
****
Peak Number 4 4-(7-Methyloctyl)phenol Concentration Rank 3

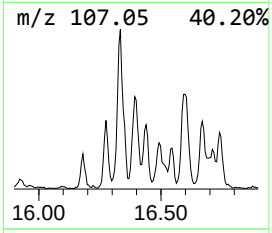
R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	6.51 ng/ul	175998	Phenanthrene-d10	17.245

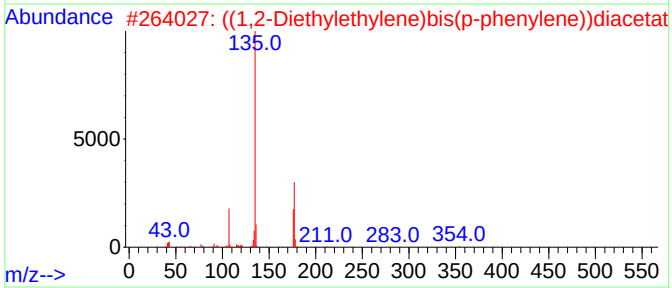
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	83
2	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	74
3	Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	64
4	Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	64
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64

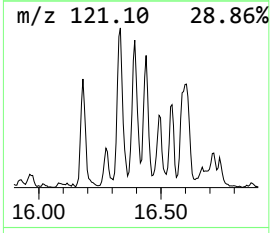


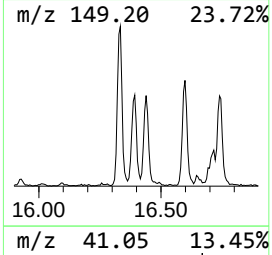


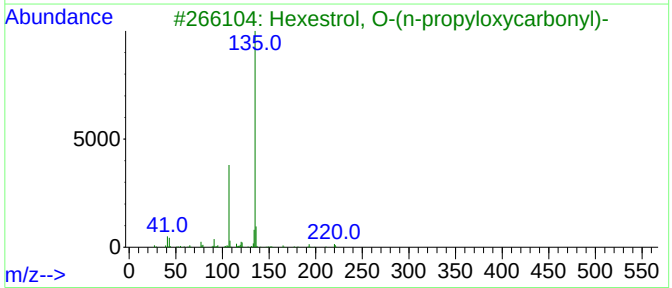


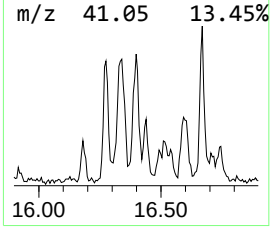












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

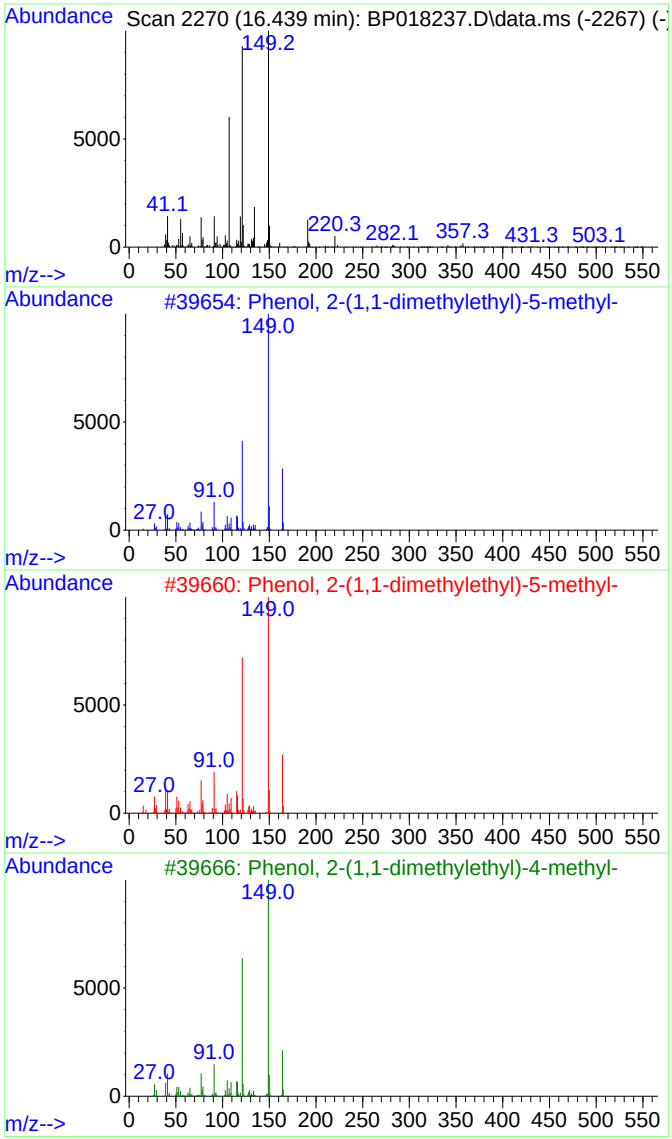
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

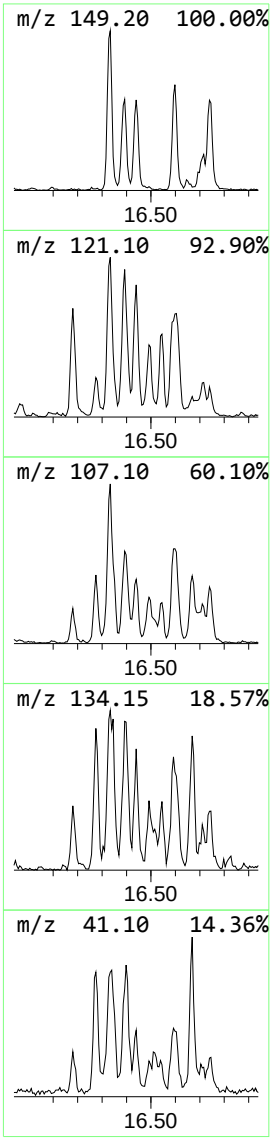
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Phenol, 2-(1,1-dimethylethy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.439	3.02 ng/ul	81787	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	53
2	Phenol, 2-(1,1-dimethylethyl)-5-...	164	C11H16O	000088-60-8	53
3	Phenol, 2-(1,1-dimethylethyl)-4-...	164	C11H16O	002409-55-4	53
4	3',4'-Formoxylidide	149	C9H11NO	006639-60-7	38
5	p-Sec-butylphenyl acetate	192	C12H16O2	003245-24-7	30





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

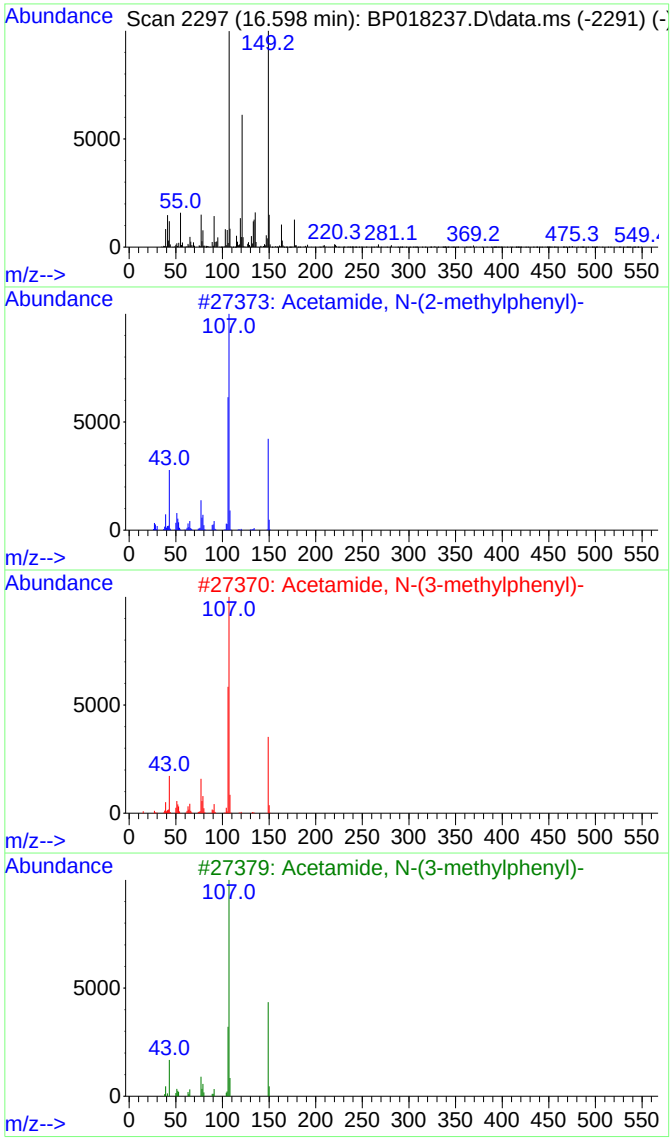
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

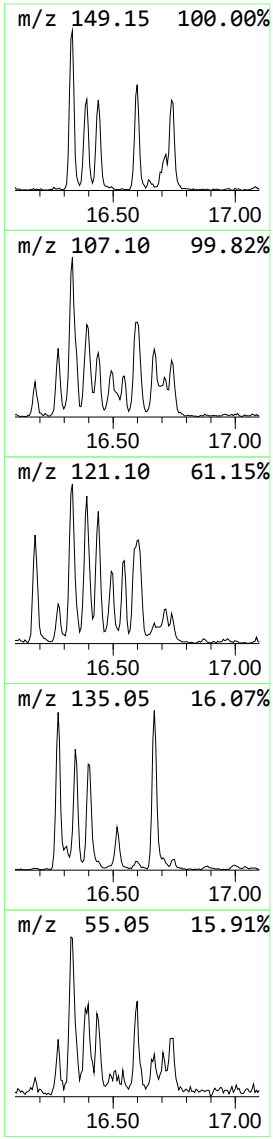
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Acetamide, N-(2-methylphenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	5.15 ng/ul	139261	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	53
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	53
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
4	Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	38
5	Benzeneacetaldehyde, 2-methoxy-....	178	C11H14O2	053155-90-1	35





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

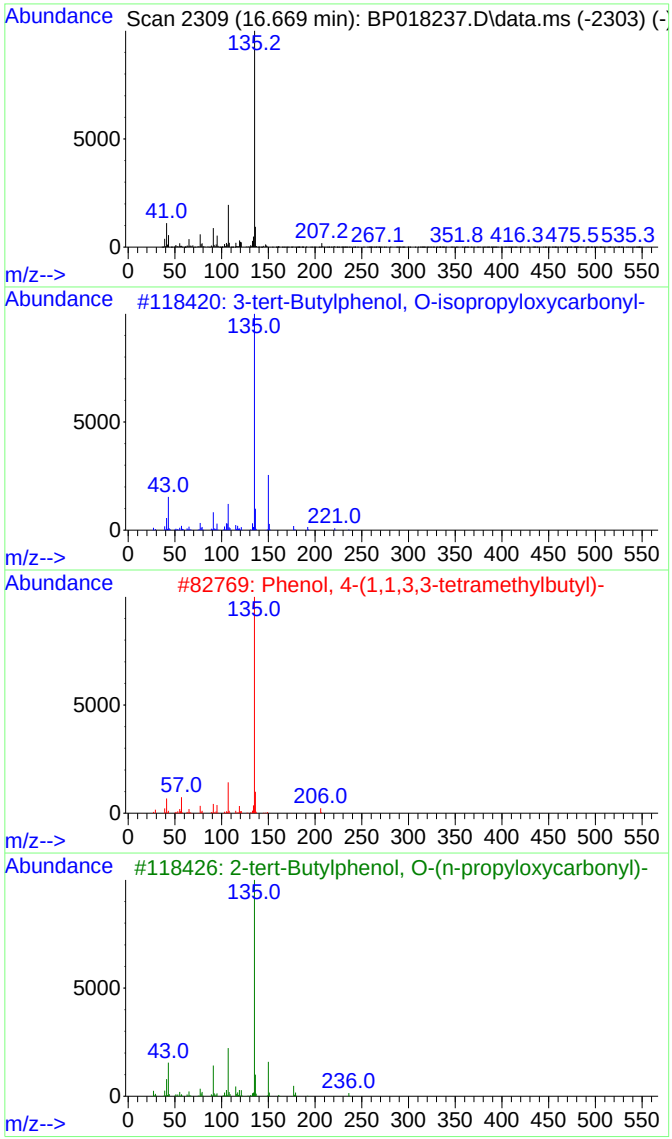
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

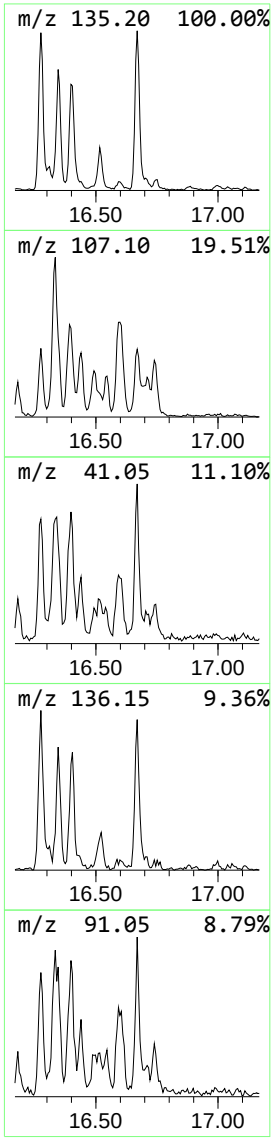
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 3-tert-Butylphenol, O-isopr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	5.55 ng/ul	150108	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-tert-Butylphenol, O-isopropylo...	236	C14H20O3	1000487-38-3	78
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
3	2-tert-Butylphenol, O-(n-propylo...	236	C14H20O3	1000487-37-9	72
4	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	72
5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

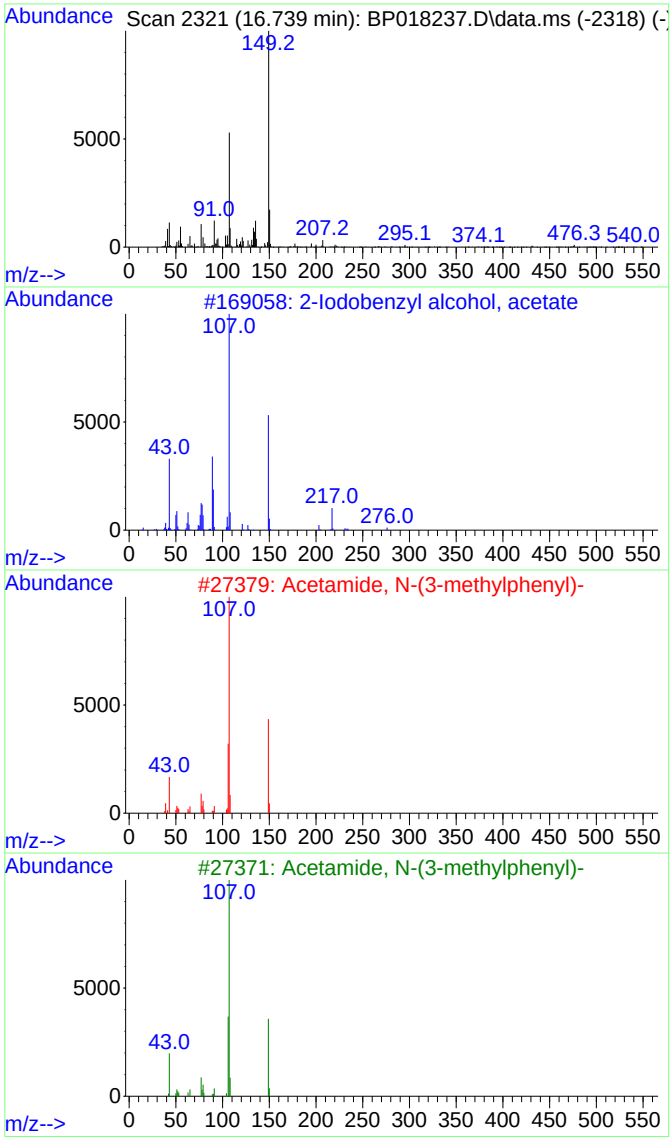
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

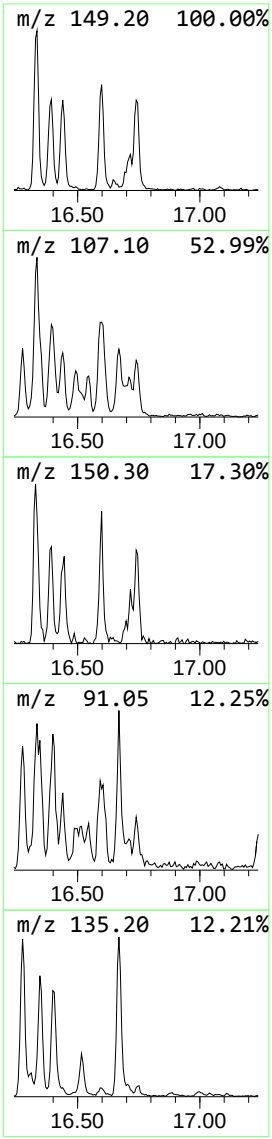
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-Iodobenzyl alcohol, acetate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.739	2.37 ng/ul	63989	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Iodobenzyl alcohol, acetate	276	C9H9IO2	1000364-47-1	64
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
4	3-Methyl-1-adamantanecarboxylic ...	194	C12H18O2	1000504-38-9	50
5	Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

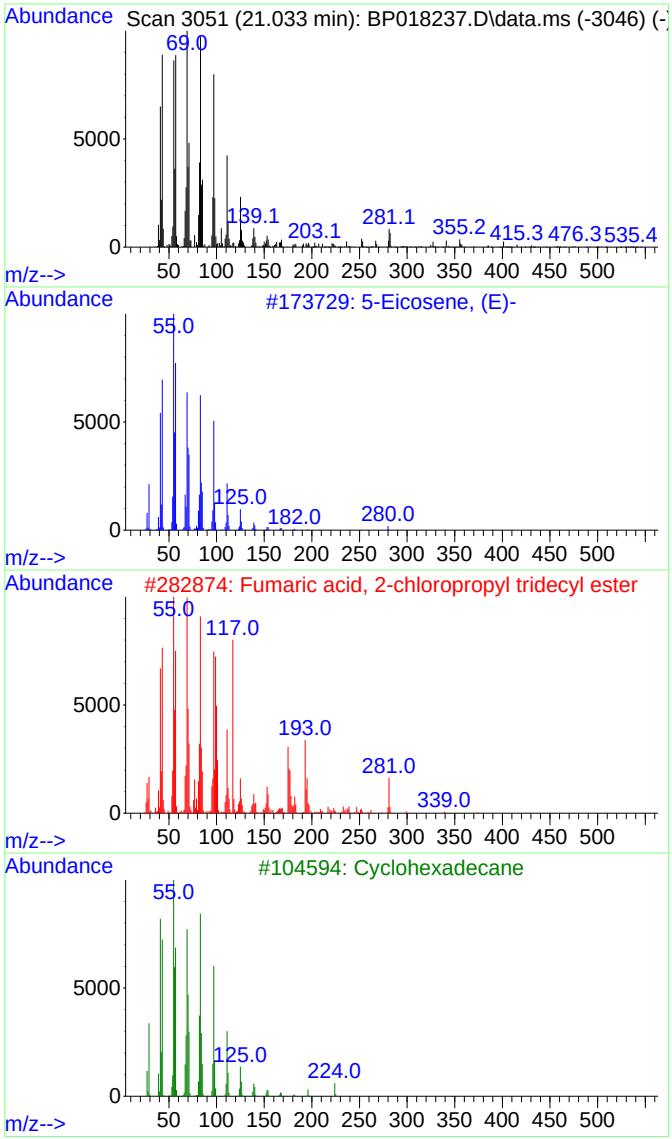
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

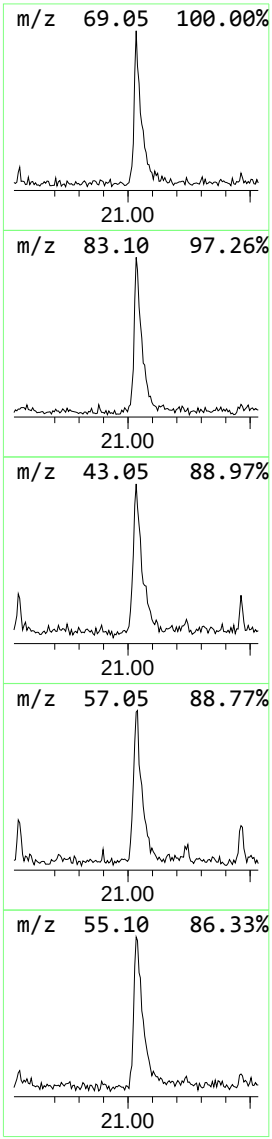
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.033	5.91 ng/ul	185388	Chrysene-d12	21.374

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	97
3	Cyclohexadecane	224	C16H32	000295-65-8	95
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	95





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018237.D
Acq On : 18 Nov 2023 00:47
Operator : MA/JU
Sample : 05369-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.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
2,4,7,9-Tetrame...	13.298	35.3	ng/ul	787451	3	14.463	445623	20.0
4-(1,1-Dimethyl...	16.275	4.8	ng/ul	129053	4	17.245	541039	20.0
unknown-01	16.333	9.2	ng/ul	248530	4	17.245	541039	20.0
4-(7-Methylocty...	16.398	6.5	ng/ul	175998	4	17.245	541039	20.0
Phenol, 2-(1,1-...	16.439	3.0	ng/ul	81787	4	17.245	541039	20.0
Acetamide, N-(2...	16.598	5.2	ng/ul	139261	4	17.245	541039	20.0
3-tert-Butylphe...	16.669	5.5	ng/ul	150108	4	17.245	541039	20.0
2-Iodobenzyl al...	16.739	2.4	ng/ul	63989	4	17.245	541039	20.0
(DEL) Alkane: S...	21.033	5.9	ng/ul	185388	5	21.374	627320	20.0