LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : GCDP5

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: BP018236.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.158	9	12	19	rVB	12144	18954	1.57%	0.159%
2	3.293	33	35	38	rVB4	2231	2140	0.18%	0.018%
3	3.623	87	91	104	rBV2	17491	40490	3.36%	0.340%
4	3.711	104	106	109	rVV4	2572	2919	0.24%	0.025%
5	3.770	113	116	124	rVB6	4295	8977	0.74%	0.075%
6	3.905	135	139	140	rVB4	2606	2393	0.20%	0.020%
7	3.923	140	142	144	rVB3	1874	1821	0.15%	0.015%
8	4.740	275	281	282	rBV5	1445	1994	0.17%	0.017%
9	5.081	337	339	343	rBV5	1361	2047	0.17%	0.017%
10	5.128	343	347	350	rVB6	1357	1632	0.14%	0.014%
11	5.164	350	353	356	rBV5	1635	1757	0.15%	0.015%
12	5.487	404	408	411	rVB6	1514	2566	0.21%	0.022%
13	5.523	411	414	417	rVB5	1171	1604	0.13%	0.013%
14	6.017	496	498	503	rBV6	1023	1732	0.14%	0.015%
15	6.464	571	574	575	rBV2	2115	1944	0.16%	0.016%
16	6.552	586	589	591	rBV4	1650	2034	0.17%	0.017%
17	6.728	616	619	623	rVB6	1965	3634	0.30%	0.031%
18	6.981	654	662	664	rBV3	12768	21784	1.81%	0.183%
19	7.134	681	688	699	rBV	118463	200205	16.61%	1.683%
20	7.311	711	718	733	rBV	91834	185901	15.43%	1.563%
21	7.781	790	798	815	rBV2	144038	275302	22.84%	2.315%
22	8.522	919	924	942	rBV	52249	143019	11.87%	1.202%
23	8.993	997	1004	1023	rBV	145266	314554	26.10%	2.645%
24	9.158	1028	1032	1036	rVB6	4122	7006	0.58%	0.059%
25	9.711	1118	1126	1147	rBV	89786	217295	18.03%	1.827%
26	9.975	1167	1171	1174	rVB6	1757	2028	0.17%	0.017%
27	10.011	1174	1177	1180	rBV5	1437	1620	0.13%	0.014%
28	10.146	1198	1200	1203	rBV4	1464	2003	0.17%	0.017%
29	10.246	1210	1217	1244	rBV	96420	277192	23.00%	2.330%
30	10.522	1263	1264	1267	rBV3	1546	1729	0.14%	0.015%
31	10.599	1267	1277	1286	rBV	210731	360115	29.88%	3.028%
32	10.787	1303	1309	1333	rBV	133337	383861	31.85%	3.227%
33	11.387	1407	1411	1412	rVV4	1859	2251	0.19%	0.019%
34	11.410	1412	1415	1421	rVV8	2725	6025	0.50%	0.051%
35	11.481	1421	1427	1430	rVV8	1176	2005	0.17%	0.017%
36	11.522	1430	1434	1438	rBV6	1243	2060	0.17%	0.017%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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	11.652	1455	1456	1460	rVB4	2018	1691	0.14%	0.014%
38	12.269	1555	1561	1562	rBV6	2673	2911	0.24%	0.024%
39	12.340	1570	1573	1577	rVB5	1129	1635	0.14%	0.014%
40	12.393	1580	1582	1586	rBV5	1211	1758	0.15%	0.015%
41	13.069	1694	1697	1700	rVB5	1491	2018	0.17%	0.017%
42	13.210	1716	1721	1725	rBV8	2268	4699	0.39%	0.040%
43	13.299	1730	1736	1749	rBV2	25735	50123	4.16%	0.421%
44	13.663	1794	1798	1800	rBV5	1865	2747	0.23%	0.023%
45	13.804	1818	1822	1824	rBV5	2001	2822	0.23%	0.024%
46	13.887	1829	1836	1850	rBV	408228	625327	51.89%	5.257%
47	14.157	1876	1882	1893	rBV	428096	650455	53.98%	5.468%
48	14.287	1902	1904	1908	rVB5	1660	1807	0.15%	0.015%
49	14.363	1915	1917	1922	rVB5	1379	1890	0.16%	0.016%
50	14.463	1927	1934	1947	rBV	301049	463091	38.43%	3.893%
51	14.893	2001	2007	2012	rBV10	4163	10228	0.85%	0.086%
52	14.969	2017	2020	2023	rVV5	2017	2911	0.24%	0.024%
53	15.104	2038	2043	2045	rBV6	1923	2524	0.21%	0.021%
54	15.246	2064	2067	2072	rVB7	2271	3376	0.28%	0.028%
55	15.310	2074	2078	2080	rBV5	1833	1945	0.16%	0.016%
56	15.463	2098	2104	2114	rBV	541984	848813	70.43%	7.136%
57	15.634	2128	2133	2144	rBV	67723	151718	12.59%	1.276%
58	15.881	2170	2175	2179	rBV2	11760	14893	1.24%	0.125%
59	15.922	2179	2182	2187	rVV3	7707	11881	0.99%	0.100%
60	15.957	2187	2188	2192	rVV4	3280	3415	0.28%	0.029%
61	16.028	2196	2200	2203	rVB6	1537	2089	0.17%	0.018%
62	16.181	2221	2226	2235	rVB	22479	31490	2.61%	0.265%
63	16.275	2235	2242	2246	rBV	65852	90300	7.49%	0.759%
64	16.334	2247	2252	2258	rVV3	81864	160613	13.33%	1.350%
65	16.398	2258	2263	2267	rVV3	63654	112499	9.34%	0.946%
66	16.440	2267	2270	2275	rVV	44024	59245	4.92%	0.498%
67	16.493	2275	2279	2281	rVV	22760	34115	2.83%	0.287%
68	16.516	2281	2283	2285	rVV	22164	26468	2.20%	0.223%
69	16.540	2285	2287	2292	rVV2	21573	31421	2.61%	0.264%
70	16.598	2292	2297	2304	rVV4	54527	108817	9.03%	0.915%
71	16.669	2304	2309	2315	rVV	61432	101146	8.39%	0.850%
72	16.740	2318	2321	2328	rVV	28170	42217	3.50%	0.355%
73	16.822	2332	2335	2340	rVB7	2253	3064	0.25%	0.026%
74	16.875	2342	2344	2346	rVV3	2212	2437	0.20%	0.020%
75	16.922	2349	2352	2354	rVB4	1940	2527	0.21%	0.021%
76	17.004	2362	2366	2369	rVB6	1975	3406	0.28%	0.029%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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	17.192	2394	2398	2401	rBV6	1179	2063	0.17%	0.017%
78	17.245	2401	2407	2417	rVV	407249	581688	48.27%	4.890%
79	17.345	2418	2424	2442	rVV	611820	954407	79.20%	8.024%
80	17.481	2444	2447	2454	rVB9	4757	10382	0.86%	0.087%
81	17.663	2476	2478	2480	rVV3	1899	1799	0.15%	0.015%
82	17.687	2480	2482	2484	rVV3	1916	1775	0.15%	0.015%
83	17.728	2486	2489	2492	rVV4	1483	1820	0.15%	0.015%
84	17.875	2507	2514	2520	rBV9	9497	16720	1.39%	0.141%
85	17.975	2529	2531	2534	rBV4	2400	1747	0.14%	0.015%
86	18.092	2546	2551	2563	rBV2	80132	135325	11.23%	1.138%
87	18.304	2583	2587	2593	rVV	23074	32658	2.71%	0.275%
88	18.587	2633	2635	2637	rBV3	1915	2006	0.17%	0.017%
89	18.751	2660	2663	2664	rBV3	2248	2401	0.20%	0.020%
90	18.781	2666	2668	2669	rBV2	1651	1683	0.14%	0.014%
91	19.181	2731	2736	2738	rBV6	7814	12096	1.00%	0.102%
92	19.204	2738	2740	2741	rBV2	5714	4661	0.39%	0.039%
93	19.339	2759	2763	2769	rBV2	54159	78958	6.55%	0.664%
94	19.598	2802	2807	2818	rBV	845548	1162447	96.46%	9.773%
95	20.057	2883	2885	2892	rVB7	9386	13540	1.12%	0.114%
96	20.551	2967	2969	2971	rBV3	10348	9647	0.80%	0.081%
97	21.039	3046	3052	3062	rBV4	61496	156015	12.95%	1.312%
98	21.375	3105	3109	3120	rBV	472621	638327	52.97%	5.367%
99	23.680	3495	3501	3516	rVB2	604100	1205103	100.00%	10.132%
100	23.851	3523	3530	3540	rBV	316863	686206	56.94%	5.769%

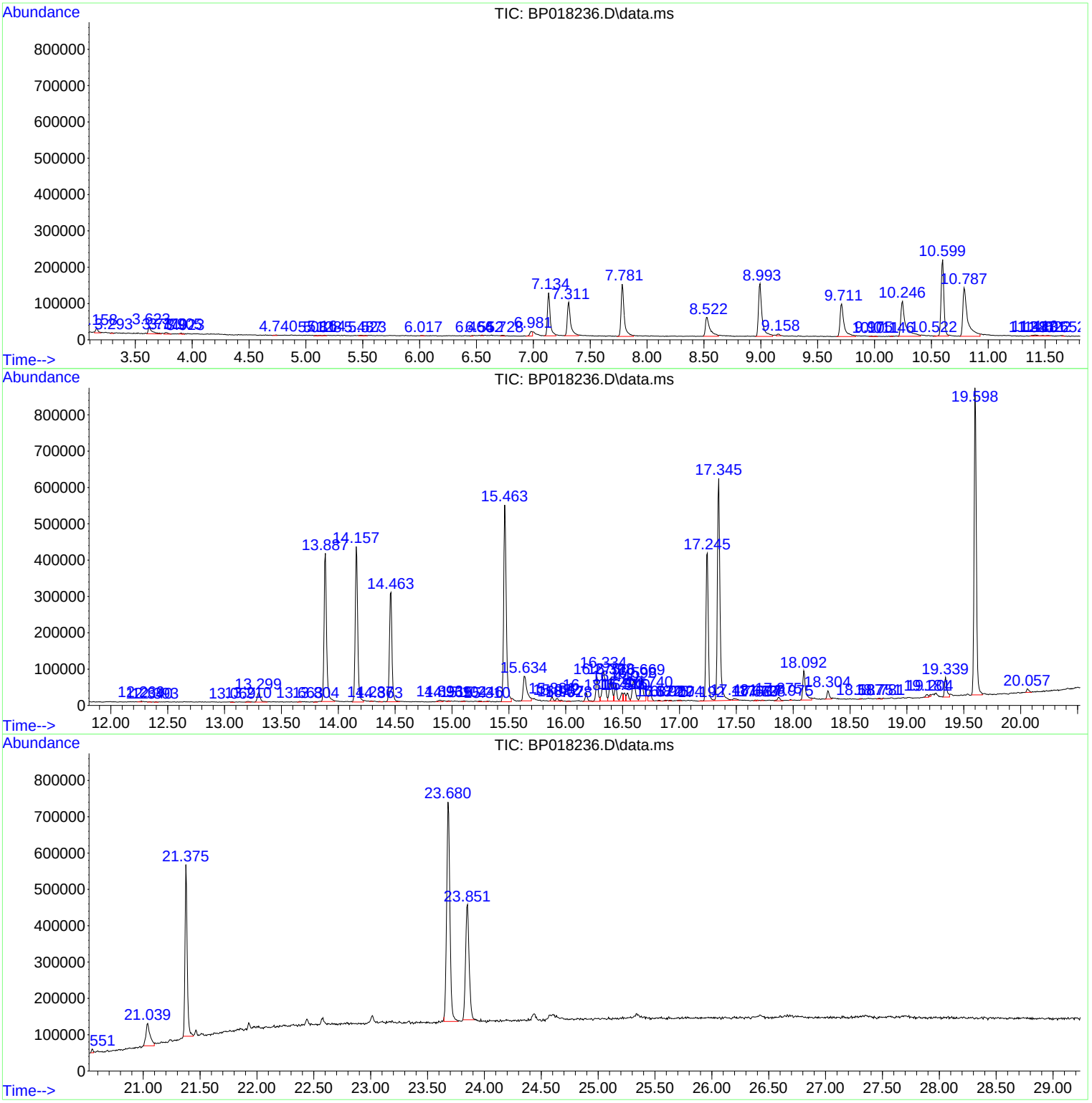
Sum of corrected areas: 11894599

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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 : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

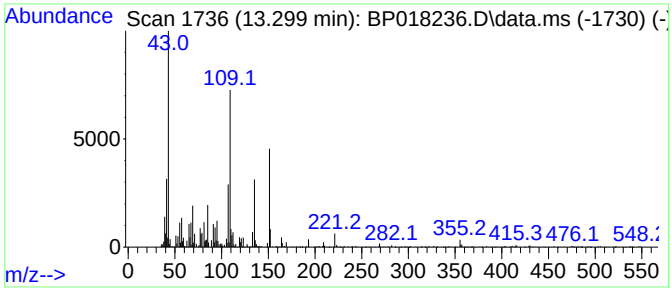
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

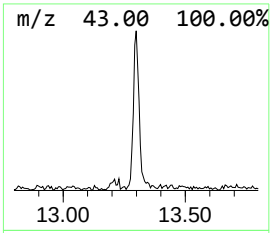
****

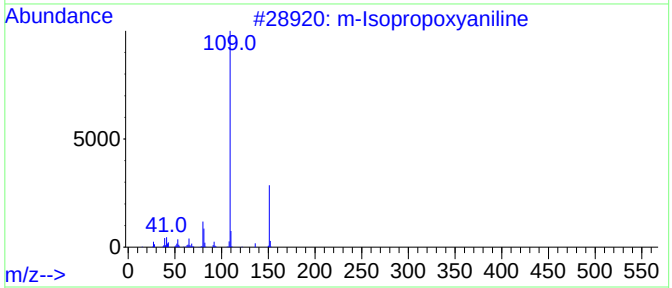
Peak Number 1 unknown-01 Concentration Rank 9

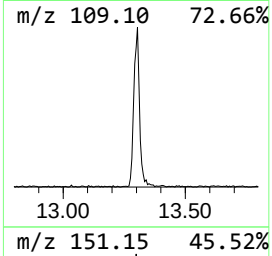
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.299	2.16 ng/ul	50123	Acenaphthene-d10	14.463

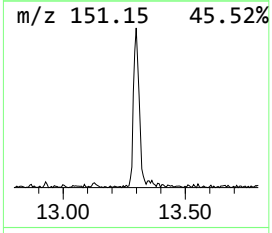
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	43
2	Metacetamol	151	C8H9NO2	000621-42-1	38
3	Butanamide, N-acetyl-N-(4-hydrox...	221	C12H15NO3	1000107-08-7	38
4	Acetaminophen	151	C8H9NO2	000103-90-2	38
5	N-Cyclopropyl-p-fluorobenzylamine	165	C10H12FN	1000250-88-9	14

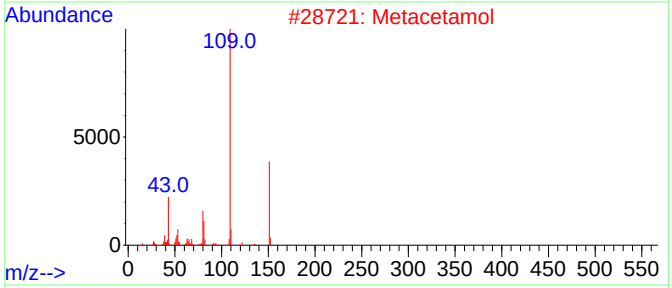


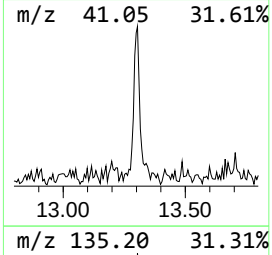


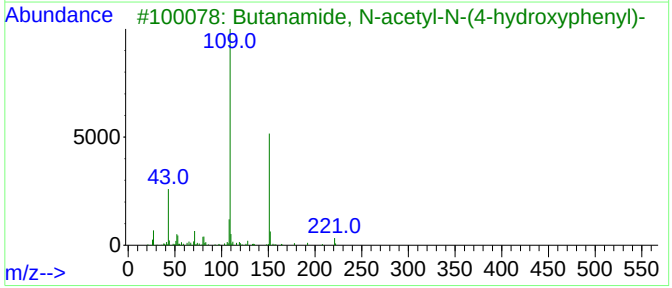


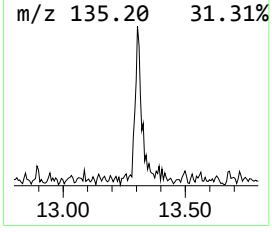












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

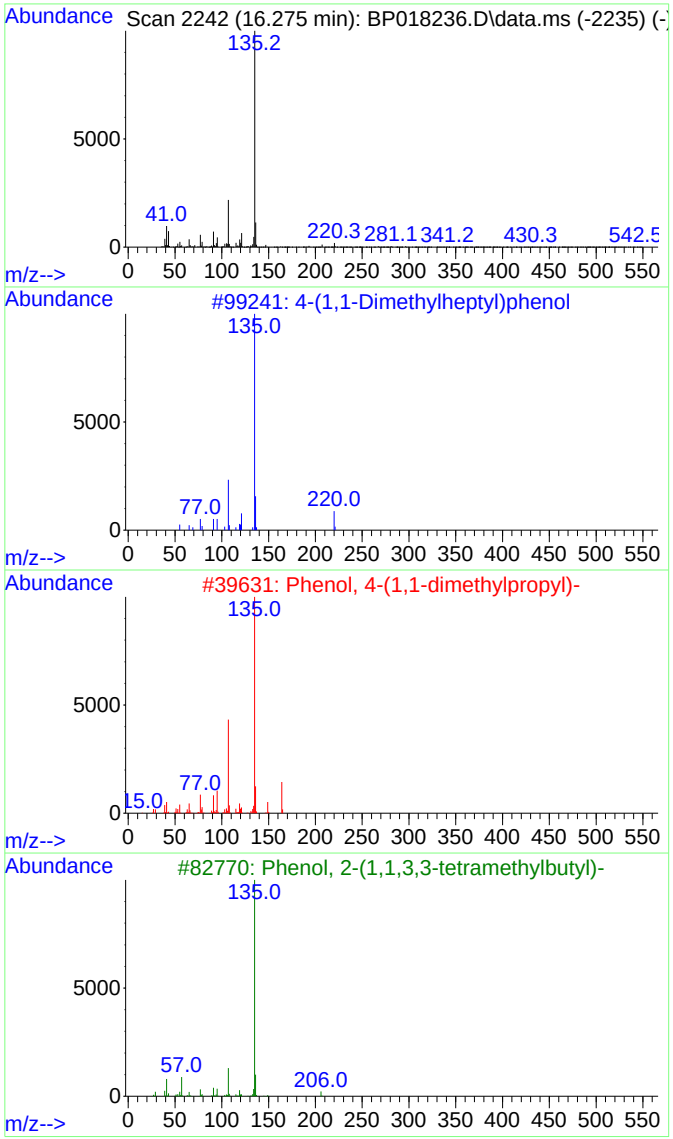
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

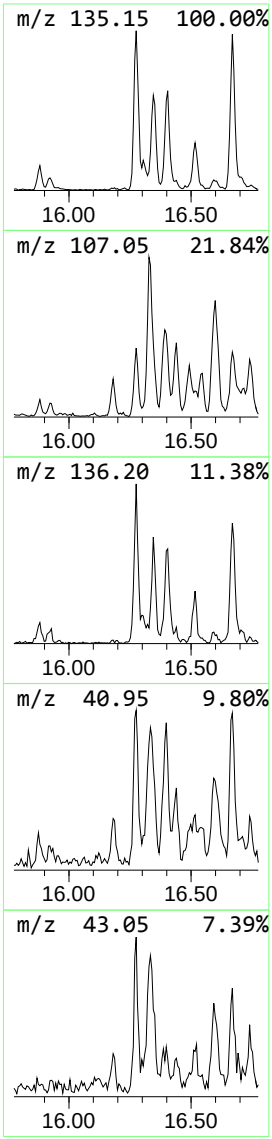
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	3.10 ng/ul	90300	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	93
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
4	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72
5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

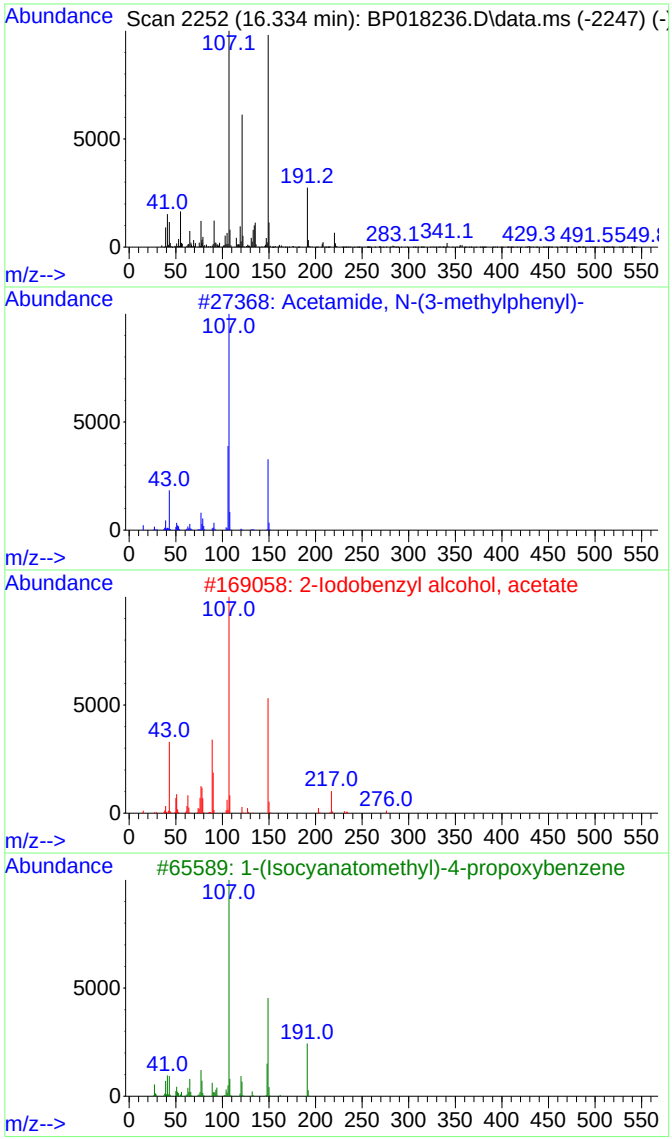
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

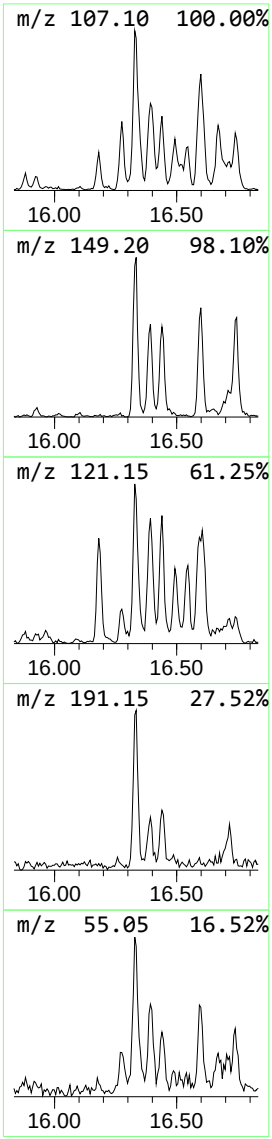
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Acetamide, N-(3-methylphenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	5.52 ng/ul	160613	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
2	2-Iodobenzyl alcohol, acetate	276	C9H9IO2	1000364-47-1	53
3	1-(Isocyanatomethyl)-4-propoxybe...	191	C11H13NO2	936095-07-7	52
4	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	46
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

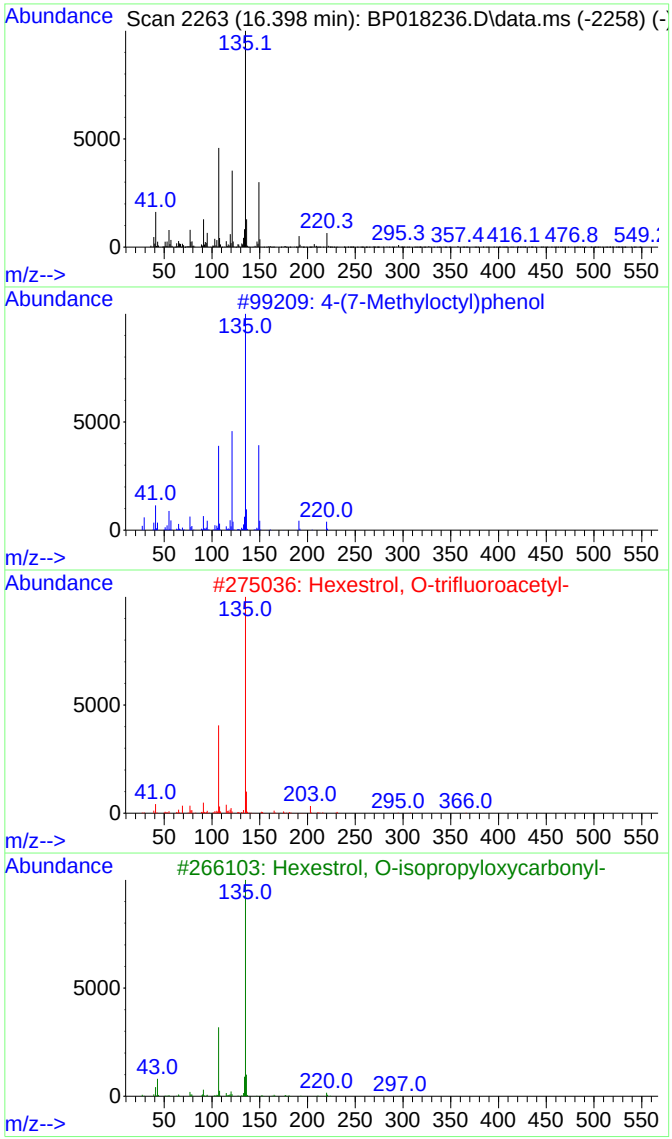
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

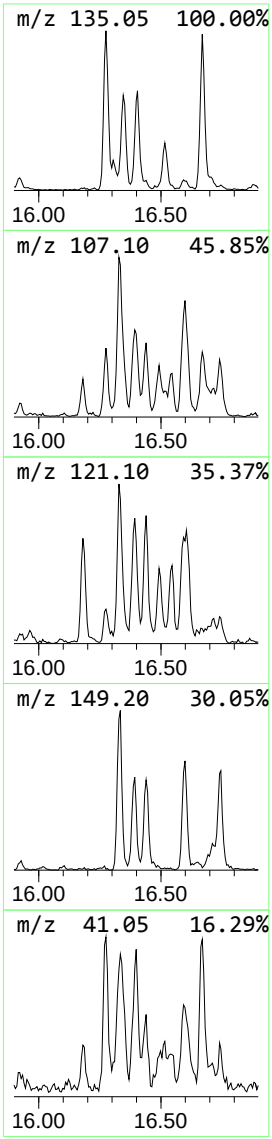
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	3.87 ng/ul	112499	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	91
2	Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	59
3	Hexestrol, O-isopropyloxycarbonyl-	356	C22H28O4	1000487-68-4	59
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	59
5	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	59





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Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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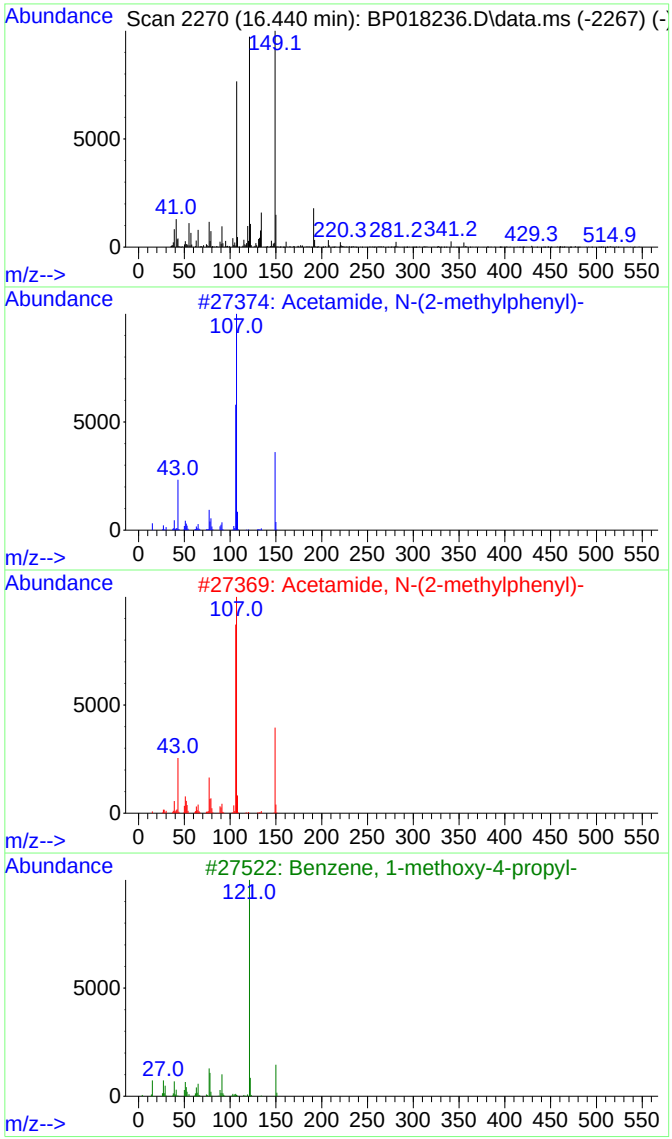
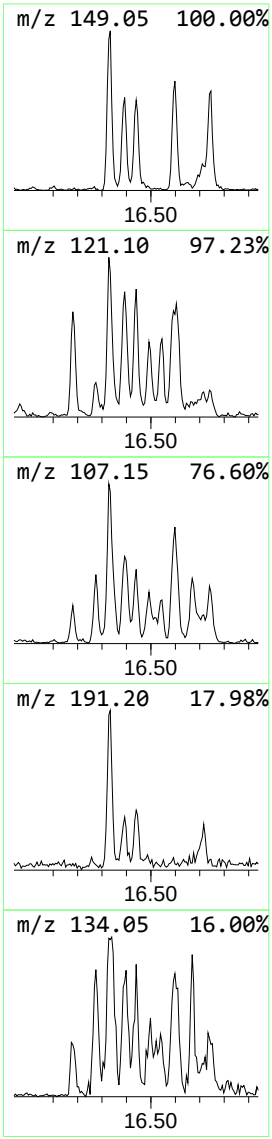
Peak Number 5 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.440	2.04 ng/ul	59245	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
3	Benzene, 1-methoxy-4-propyl-	150	C10H14O	000104-45-0	35
4	2-Methyl-6-propylphenol	150	C10H14O	003520-52-3	25
5	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	25





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Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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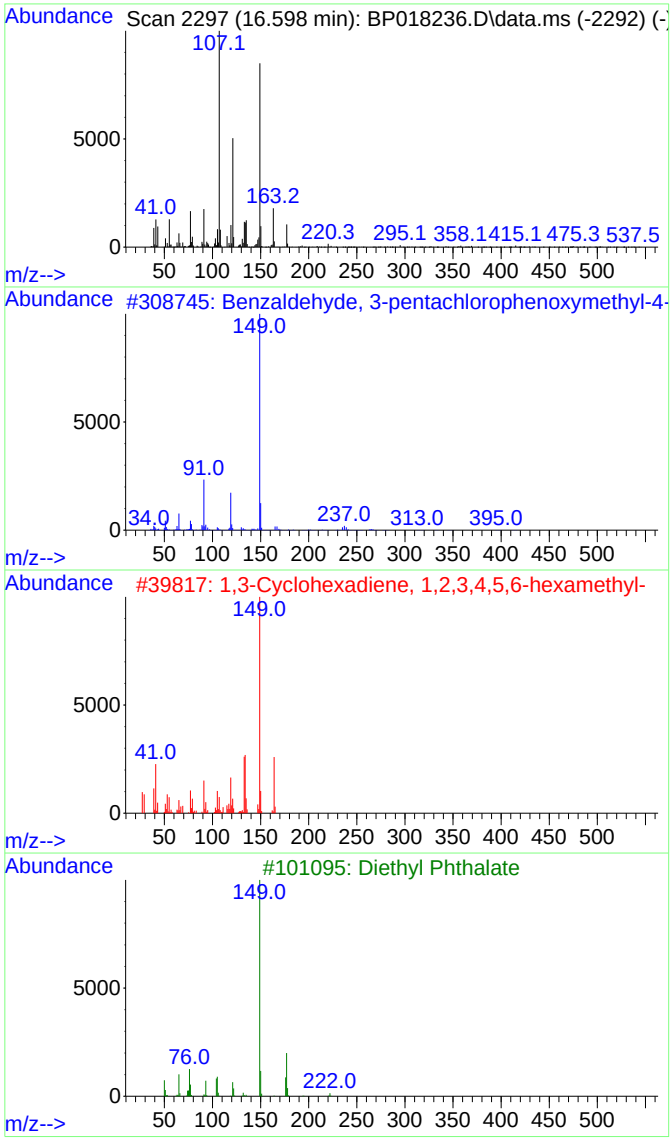
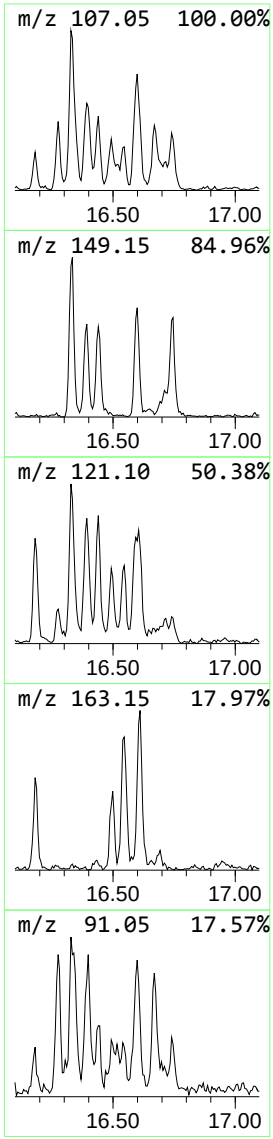
Peak Number 6 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	3.74 ng/ul	108817	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3-pentachloropheno...	412	C15H9Cl5O3	329222-74-4	38
2	1,3-Cyclohexadiene, 1,2,3,4,5,6-...	164	C12H20	1000152-09-6	38
3	Diethyl Phthalate	222	C12H14O4	000084-66-2	35
4	Phthalic acid, 7-bromoheptyl eth...	370	C17H23BrO4	1000415-51-6	35
5	Phenol, 2-(1,1-dimethylethyl)-4-...	164	C11H16O	002409-55-4	35





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Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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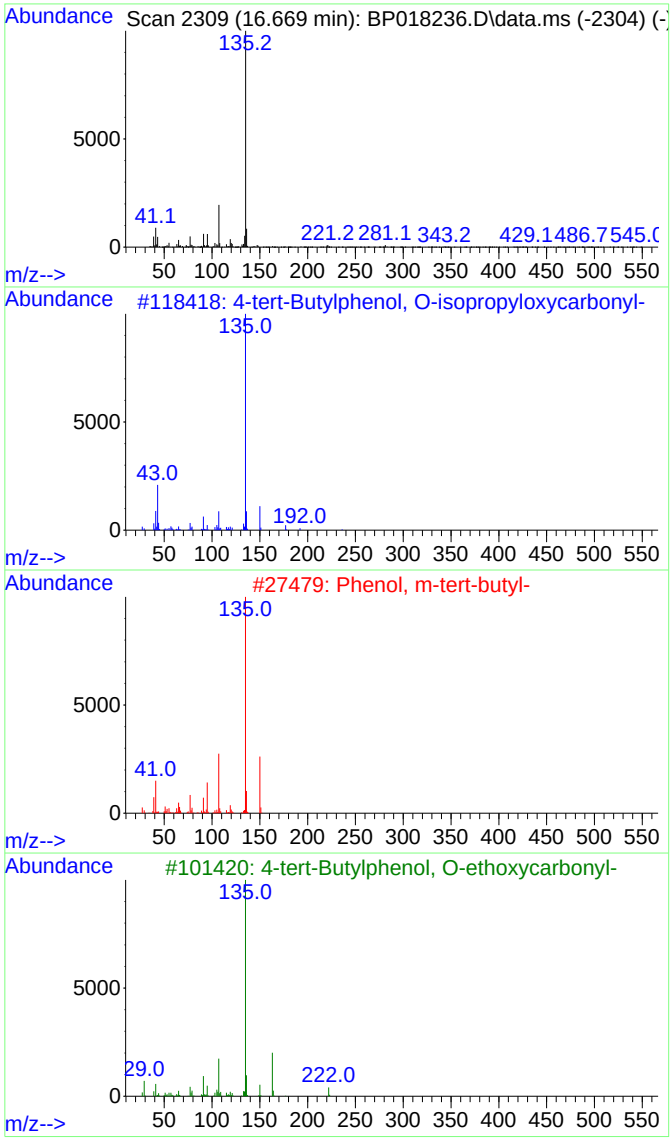
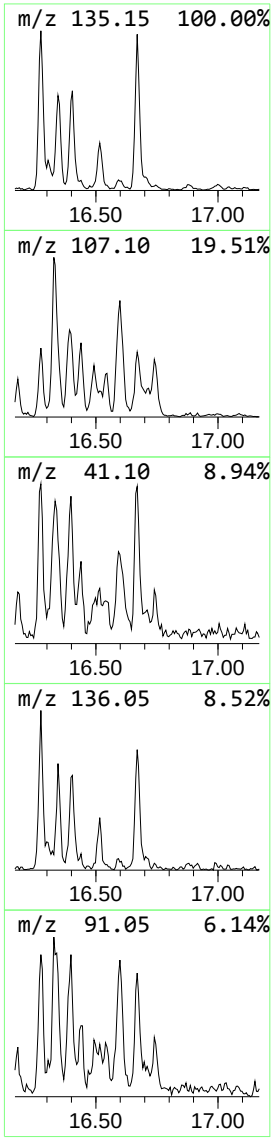
Peak Number 7 4-tert-Butylphenol, O-isopr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	3.48 ng/ul	101146	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-tert-Butylphenol, O-isopropylo...	236	C14H20O3	1201410-82-3	78
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
3	4-tert-Butylphenol, O-ethoxycarb...	222	C13H18O3	026177-66-2	78
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
5	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	72





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Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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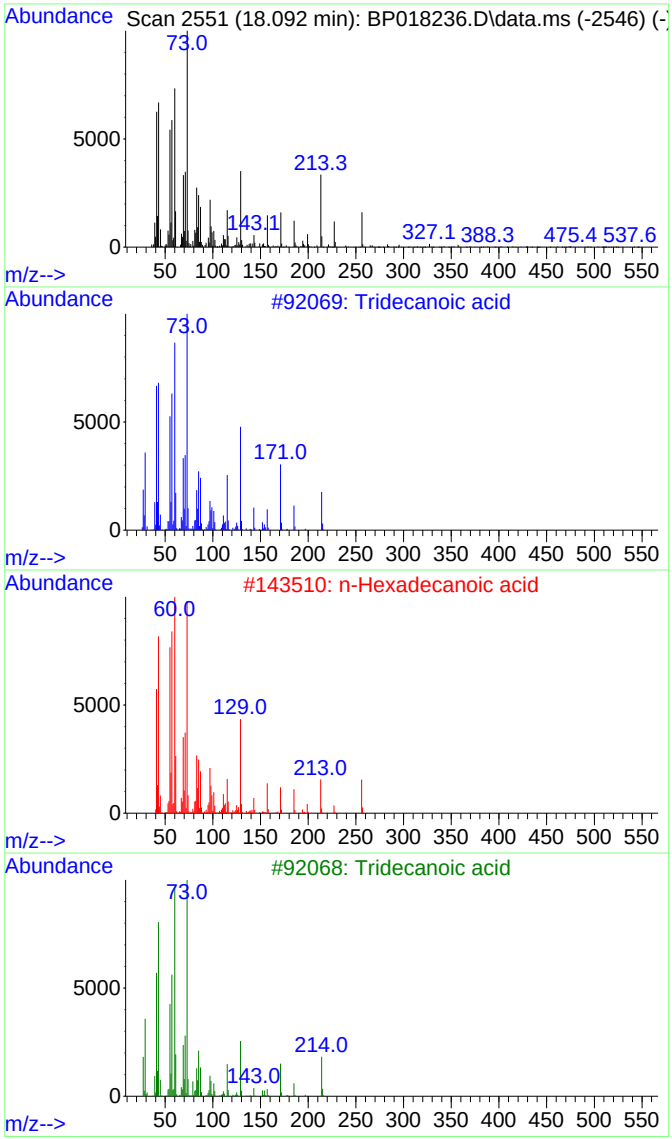
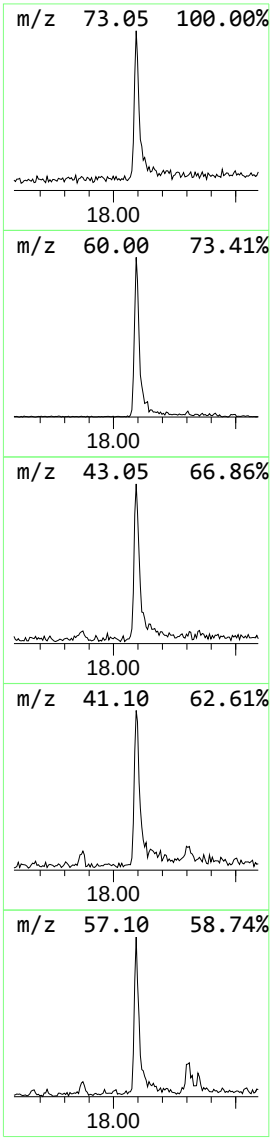
Peak Number 8 Tridecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	4.65 ng/ul	135325	Phenanthrene-d10	17.245

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	96
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	Tridecanoic acid	214	C13H26O2	000638-53-9	95
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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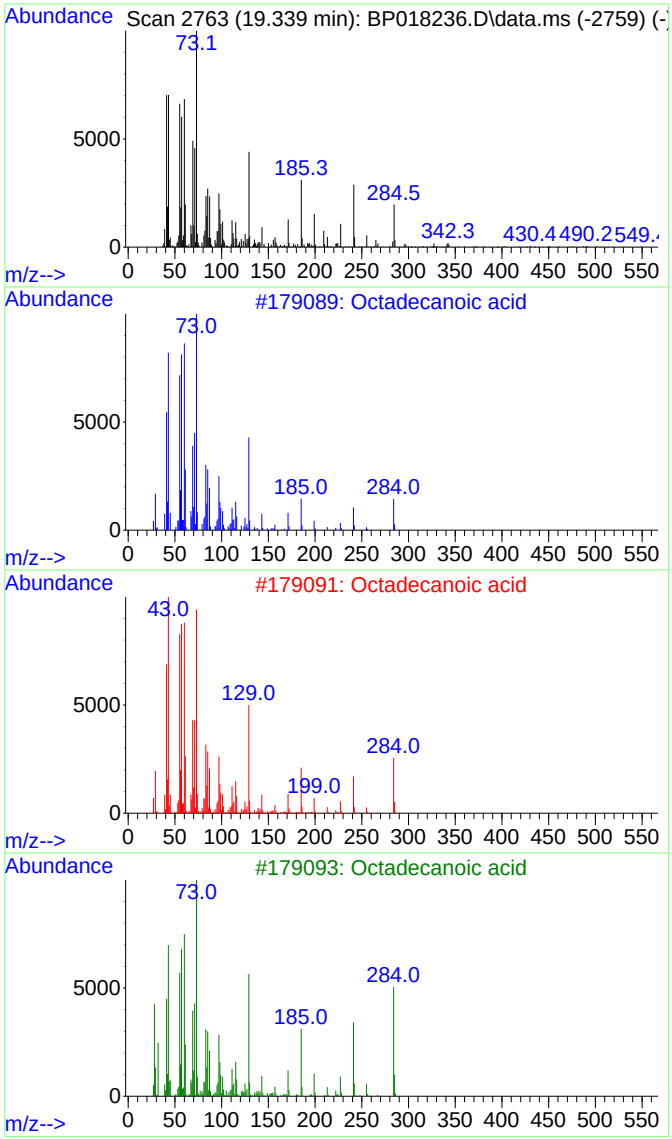
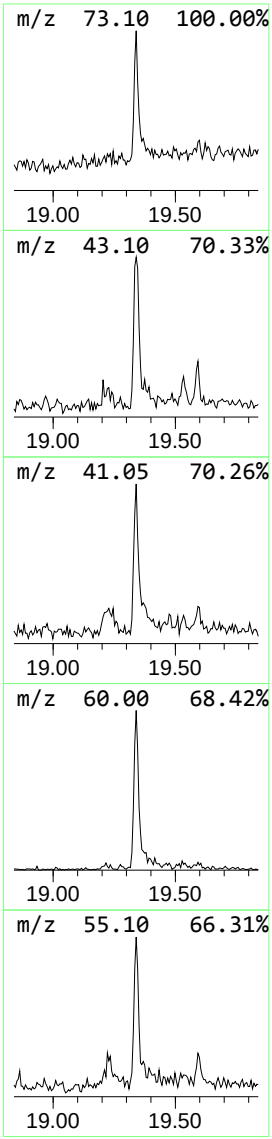
Peak Number 9 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.339	2.47 ng/ul	78958	Chrysene-d12	21.375

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Octadecanoic acid	284	C18H36O2	000057-11-4	97
3	Octadecanoic acid	284	C18H36O2	000057-11-4	96
4	Octadecanoic acid	284	C18H36O2	000057-11-4	90
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	89





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Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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

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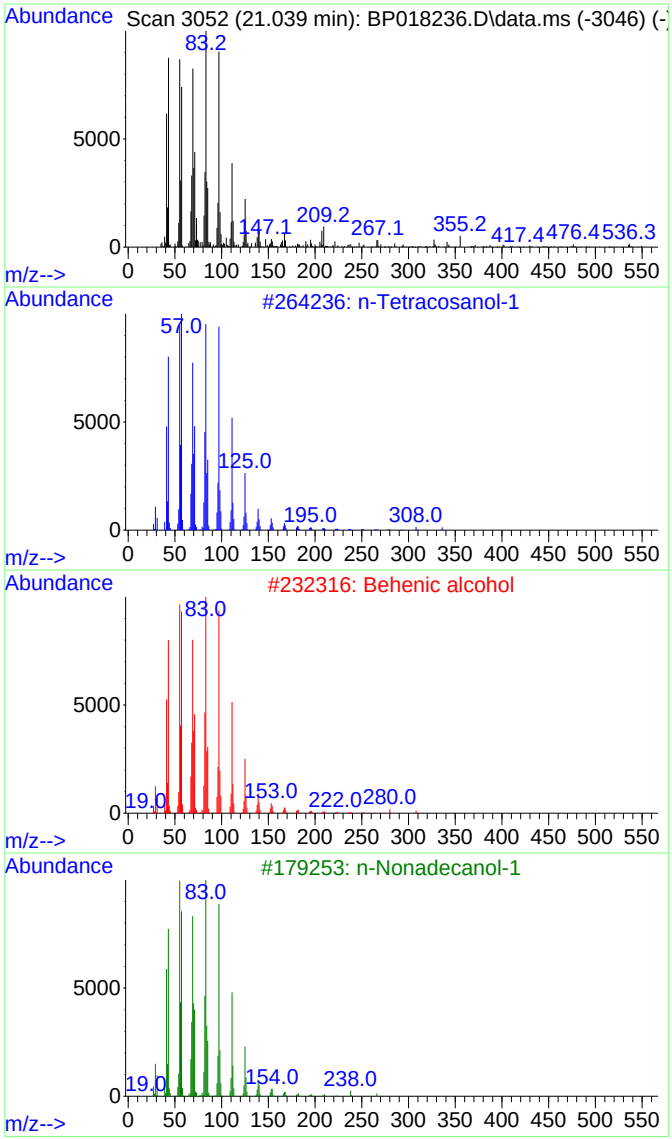
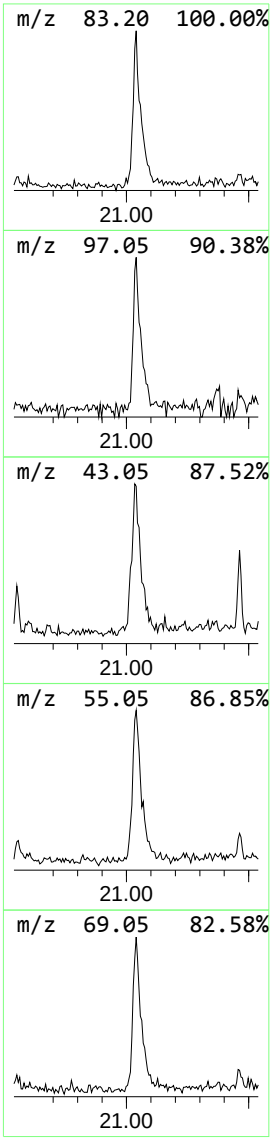
Peak Number 10 n-Tetracosanol-1 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.039	4.89 ng/ul	156015	Chrysene-d12	21.375

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
2	Behenic alcohol	326	C22H46O	000661-19-8	95
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4	1-Hexacosene	364	C26H52	018835-33-1	91
5	1-Heneicosanol	312	C21H44O	015594-90-8	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018236.D
Acq On : 18 Nov 2023 00:13
Operator : MA/JU
Sample : 05369-20
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDP5

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
unknown-01	13.299	2.2	ng/ul	50123	3	14.463	463091	20.0
4-(1,1-Dimethyl...	16.275	3.1	ng/ul	90300	4	17.245	581688	20.0
Acetamide, N-(3...	16.334	5.5	ng/ul	160613	4	17.245	581688	20.0
4-(7-Methylocty...	16.398	3.9	ng/ul	112499	4	17.245	581688	20.0
unknown-02	16.440	2.0	ng/ul	59245	4	17.245	581688	20.0
unknown-03	16.598	3.7	ng/ul	108817	4	17.245	581688	20.0
4-tert-Butylphe...	16.669	3.5	ng/ul	101146	4	17.245	581688	20.0
Tridecanoic acid	18.092	4.7	ng/ul	135325	4	17.245	581688	20.0
Octadecanoic acid	19.339	2.5	ng/ul	78958	5	21.375	638327	20.0
n-Tetracosanol-1	21.039	4.9	ng/ul	156015	5	21.375	638327	20.0