LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCFN2

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-BP060723.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP015680.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	44	rVB	49704	64998	0.77%	0.112%
2	3.846	109	112	115	rVB2	8865	9446	0.11%	0.016%
3	3.899	118	121	124	rBV5	6682	9606	0.11%	0.017%
4	3.946	125	129	136	rVB	46761	63987	0.76%	0.110%
5	4.023	138	142	145	rBV5	6260	9514	0.11%	0.016%
6	4.070	146	150	154	rVB	20017	28124	0.33%	0.049%
7	4.175	164	168	173	rVB3	13502	21093	0.25%	0.036%
8	4.411	204	208	213	rVB2	18290	24616	0.29%	0.043%
9	4.552	229	232	238	rVB3	25652	37620	0.44%	0.065%
10	4.752	263	266	270	rVB5	8298	11296	0.13%	0.020%
11	5.117	325	328	333	rBV3	5635	9332	0.11%	0.016%
12	5.240	345	349	354	rVB4	8523	15710	0.19%	0.027%
13	5.675	417	423	427	rBV9	3977	8554	0.10%	0.015%
14	6.046	480	486	490	rVB7	9304	18387	0.22%	0.032%
15	6.111	492	497	501	rVB8	5217	8902	0.11%	0.015%
16	6.305	526	530	534	rBV6	5313	9445	0.11%	0.016%
17	7.099	662	665	671	rBV8	5986	10333	0.12%	0.018%
18	7.246	681	690	703	rBV	617419	1206659	14.25%	2.084%
19	7.399	710	716	736	rVB	552025	950608	11.23%	1.642%
20	7.605	745	751	761	rBV	746732	1246211	14.72%	2.152%
21	8.081	825	832	844	rBV	560093	942273	11.13%	1.627%
22	8.716	933	940	946	rBV	4497	11397	0.13%	0.020%
23	8.805	946	955	969	rBV	583368	1099109	12.98%	1.898%
24	8.922	969	975	983	rVB	90375	166265	1.96%	0.287%
25	9.258	1025	1032	1046	rBV	657526	1213667	14.33%	2.096%
26	9.410	1053	1058	1065	rVB9	12986	27270	0.32%	0.047%
27	9.628	1090	1095	1101	rVB5	6741	12202	0.14%	0.021%
28	9.987	1146	1156	1167	rBV	585284	1054227	12.45%	1.821%
29	10.205	1186	1193	1200	rBV2	23259	72031	0.85%	0.124%
30	10.352	1214	1218	1226	rVB3	20135	37456	0.44%	0.065%
31	10.540	1242	1250	1272	rBV	618688	1360292	16.06%	2.349%
32	10.852	1298	1303	1305	rBV4	8502	11290	0.13%	0.019%
33	10.905	1305	1312	1319	rBV	707495	1205012	14.23%	2.081%
34	11.063	1331	1339	1356	rBV	244595	616408	7.28%	1.064%
35	11.752	1447	1456	1460	rBV	7422	21057	0.25%	0.036%
36	11.869	1472	1476	1479	rBV6	5387	10491	0.12%	0.018%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DCFN2

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-BP060723.MA.M
Title : SVOA CALIBRATION

37	11.904	1479	1482	1486	rVB4	9916	14343	0.17%	0.025%
38	12.304	1544	1550	1556	rBV6	10819	22901	0.27%	0.040%
39	12.499	1579	1583	1589	rVB4	11094	20219	0.24%	0.035%
40	12.569	1589	1595	1600	rBV3	9731	18172	0.21%	0.031%
41	12.716	1612	1620	1624	rBV4	5511	12894	0.15%	0.022%
42	12.787	1626	1632	1636	rVB3	11763	21507	0.25%	0.037%
43	13.193	1693	1701	1705	rBV10	7930	19034	0.22%	0.033%
44	13.240	1705	1709	1713	rVB6	9061	12721	0.15%	0.022%
45	13.328	1718	1724	1730	rVB10	7888	15109	0.18%	0.026%
46	13.528	1754	1758	1763	rBV2	14513	20011	0.24%	0.035%
47	13.581	1763	1767	1768	rVV4	5563	8530	0.10%	0.015%
48	13.604	1770	1771	1779	rVB8	8653	10128	0.12%	0.017%
49	13.898	1815	1821	1823	rBV7	5074	10334	0.12%	0.018%
50	14.046	1839	1846	1852	rBV5	12744	28286	0.33%	0.049%
51	14.134	1852	1861	1867	rBV	1310907	1810044	21.37%	3.126%
52	14.234	1874	1878	1883	rVB8	6654	11879	0.14%	0.021%
53	14.422	1903	1910	1927	rBV	1399626	2186010	25.81%	3.775%
54	14.598	1936	1940	1946	rVB2	17465	24992	0.30%	0.043%
55	14.728	1948	1962	1971	rBV	912314	1396083	16.49%	2.411%
56	14.787	1971	1972	1980	rVB5	10709	16027	0.19%	0.028%
57	14.969	1997	2003	2021	rBV2	189955	580918	6.86%	1.003%
58	15.134	2025	2031	2041	rVB8	27723	76802	0.91%	0.133%
59	15.498	2091	2093	2098	rVV6	7264	10093	0.12%	0.017%
60	15.545	2098	2101	2106	rVB3	19833	27955	0.33%	0.048%
61	15.722	2125	2131	2148	rBV	1563871	2487872	29.38%	4.296%
62	15.869	2150	2156	2167	rVV	363637	641240	7.57%	1.107%
63	15.957	2167	2171	2181	rVB10	22106	53438	0.63%	0.092%
64	16.092	2188	2194	2206	rBV	134851	215992	2.55%	0.373%
65	16.440	2245	2253	2263	rBV6	45840	127514	1.51%	0.220%
66	16.651	2284	2289	2294	rVB3	20656	30135	0.36%	0.052%
67	16.869	2323	2326	2332	rVB5	18474	25492	0.30%	0.044%
68	17.022	2347	2352	2357	rBV4	18440	31727	0.37%	0.055%
69	17.216	2381	2385	2390	rBV5	23552	34381	0.41%	0.059%
70	17.363	2406	2410	2418	rBV8	29355	57765	0.68%	0.100%
71	17.434	2418	2422	2426	rBV7	20108	29938	0.35%	0.052%
72	17.492	2426	2432	2437	rVV	1033799	1517011	17.91%	2.620%
73	17.534	2437	2439	2443	rVV	95298	126343	1.49%	0.218%
74	17.592	2443	2449	2460	rVB	1530985	2304322	27.21%	3.979%
75	17.951	2507	2510	2514	rBV2	25093	34384	0.41%	0.059%
76	18.128	2537	2540	2546	rVB3	24817	32639	0.39%	0.056%

LSC Area Percent Report

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Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Title : SVOA CALIBRATION

77	18.239	2555	2559	2564	rBV3	51365	73191	0.86%	0.126%
78	18.292	2564	2568	2573	rBV2	49527	70616	0.83%	0.122%
79	18.363	2573	2580	2597	rBV	5647558	8468475	100.00%	14.624%
80	18.622	2621	2624	2627	rBV3	28481	40836	0.48%	0.071%
81	18.986	2684	2686	2693	rBV8	43570	76334	0.90%	0.132%
82	19.469	2759	2768	2782	rBV2	2407909	4483438	52.94%	7.742%
83	19.581	2782	2787	2807	rVV	2200994	3228462	38.12%	5.575%
84	19.822	2823	2828	2838	rVB	2184957	3275750	38.68%	5.657%
85	20.622	2962	2964	2970	rBV4	100283	171363	2.02%	0.296%
86	21.251	3067	3071	3076	rVB2	474890	687639	8.12%	1.187%
87	21.580	3122	3127	3132	rBV	1185383	1712233	20.22%	2.957%
88	22.174	3225	3228	3231	rBV	226378	258713	3.06%	0.447%
89	23.292	3413	3418	3424	rVB	842475	1400833	16.54%	2.419%
90	23.363	3427	3430	3443	rVB3	440718	1033285	12.20%	1.784%
91	23.974	3528	3534	3541	rVB2	1537268	3148379	37.18%	5.437%
92	24.145	3557	3563	3571	rVB	853842	1784450	21.07%	3.082%
93	24.745	3659	3665	3677	rVB	1028815	2243298	26.49%	3.874%

Sum of corrected areas: 57907368

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
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Operator : MA/JU
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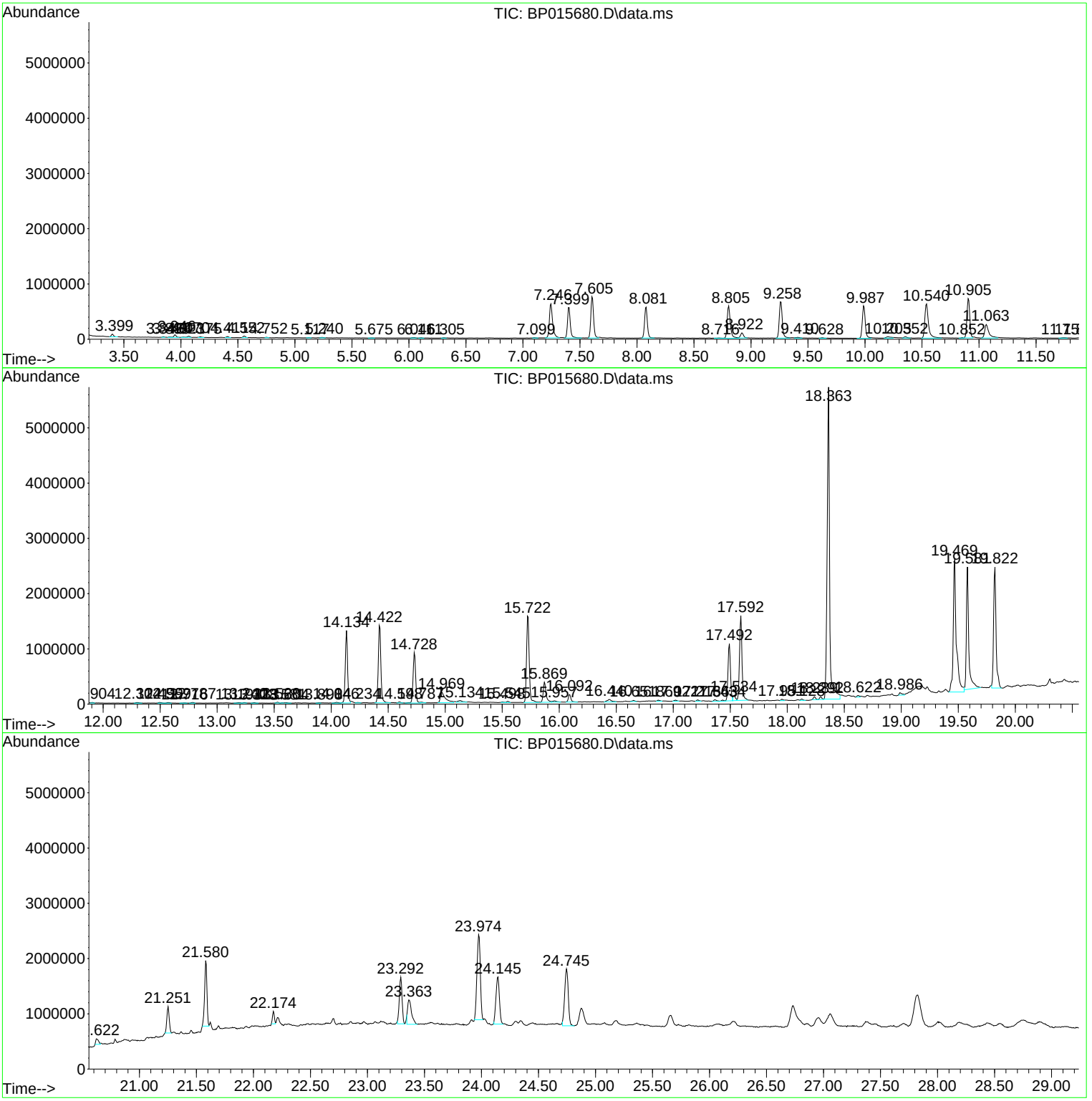
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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\BP062323\
Data File : BP015680.D
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Quant Title : SVOA CALIBRATION

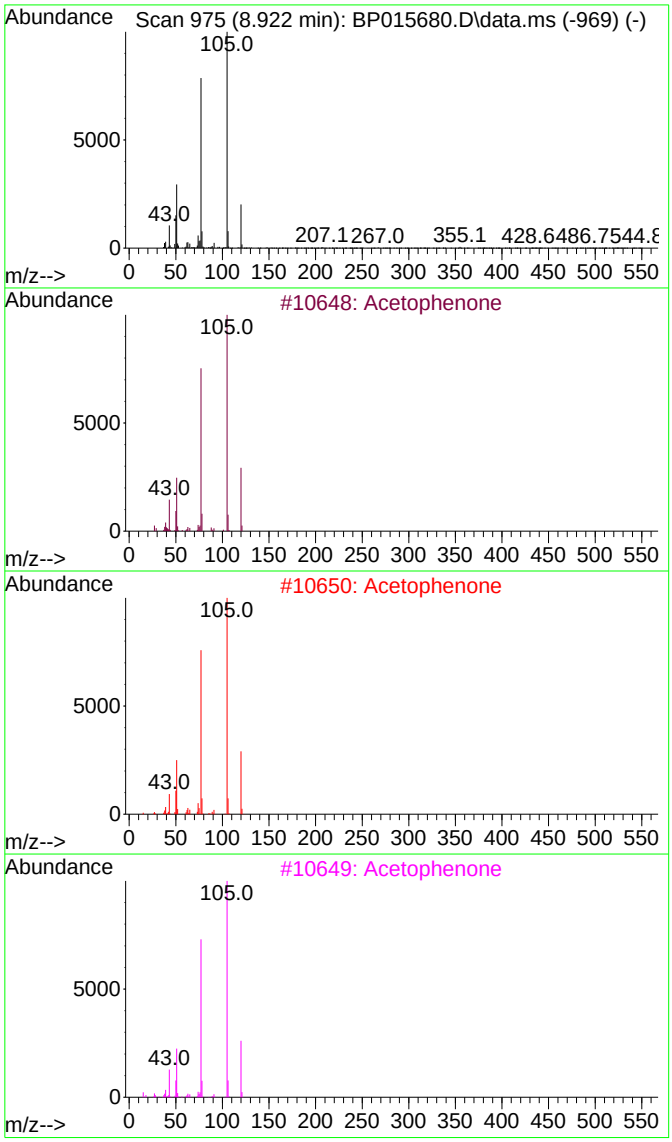
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

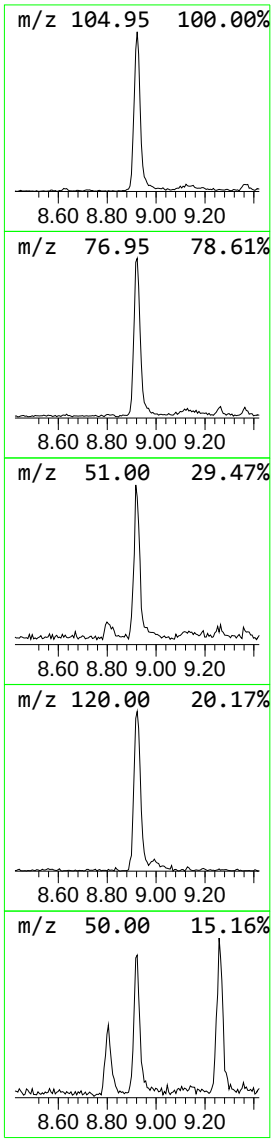
****

Peak Number 1 Aceto phenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	3.53 ng/ul	166265	1,4-Dichlorobenzene-d4	8.081

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone	120	C8H8O	000098-86-2	91
2	Acetophenone	120	C8H8O	000098-86-2	91
3	Acetophenone	120	C8H8O	000098-86-2	90
4	Acetophenone	120	C8H8O	000098-86-2	83
5	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

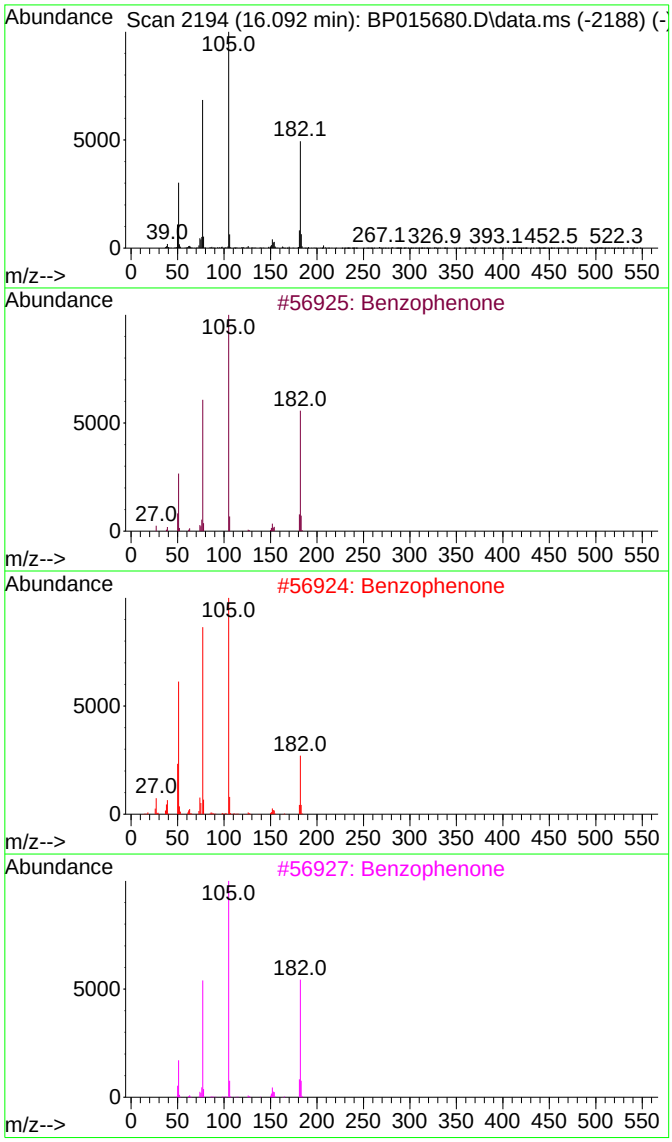
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

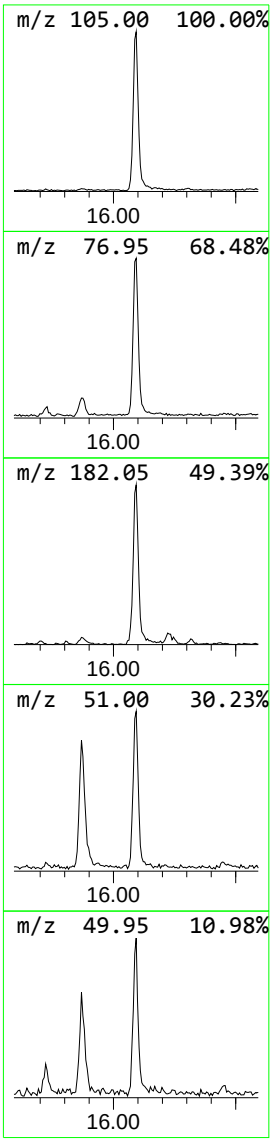
****

Peak Number 2 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.092	3.09 ng/ul	215992	Acenaphthene-d10	14.728

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	94
5	Benzophenone	182	C13H10O	000119-61-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

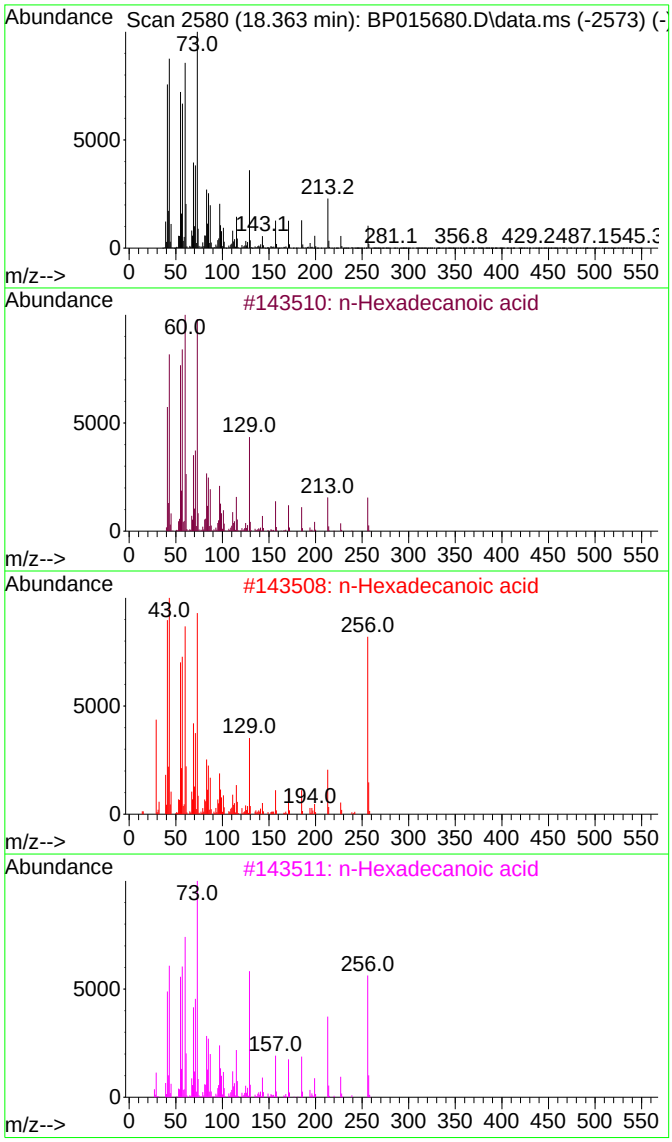
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

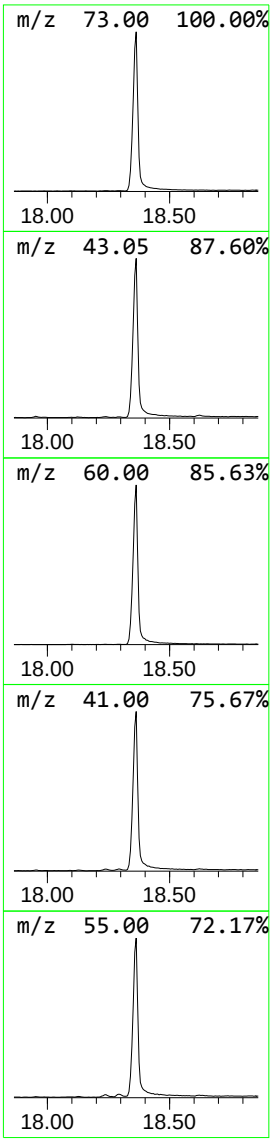
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.363	111.65 ng/ul	8468480	Phenanthrene-d10	17.492

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	97
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

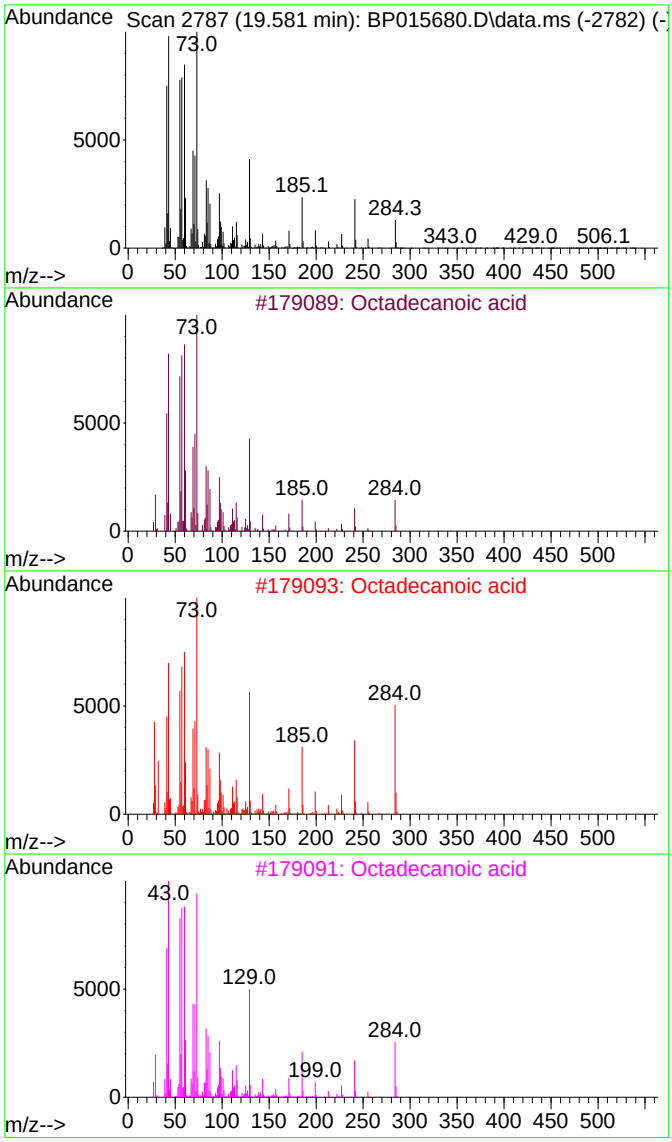
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

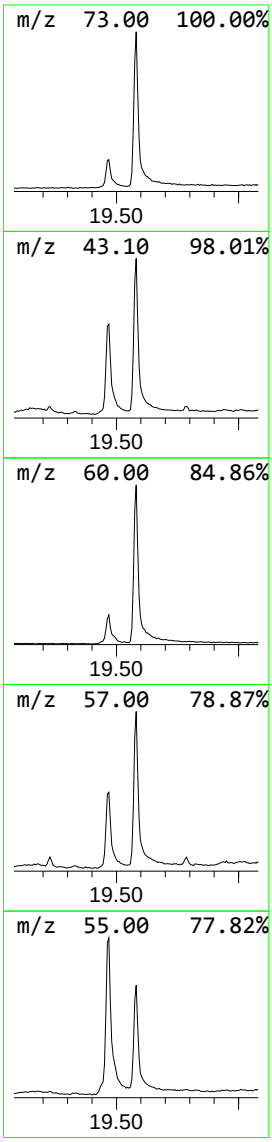
****

Peak Number 4 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.581	37.71 ng/ul	3228460	Chrysene-d12	21.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	99
4	Octadecanoic acid	284	C18H36O2	000057-11-4	95
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

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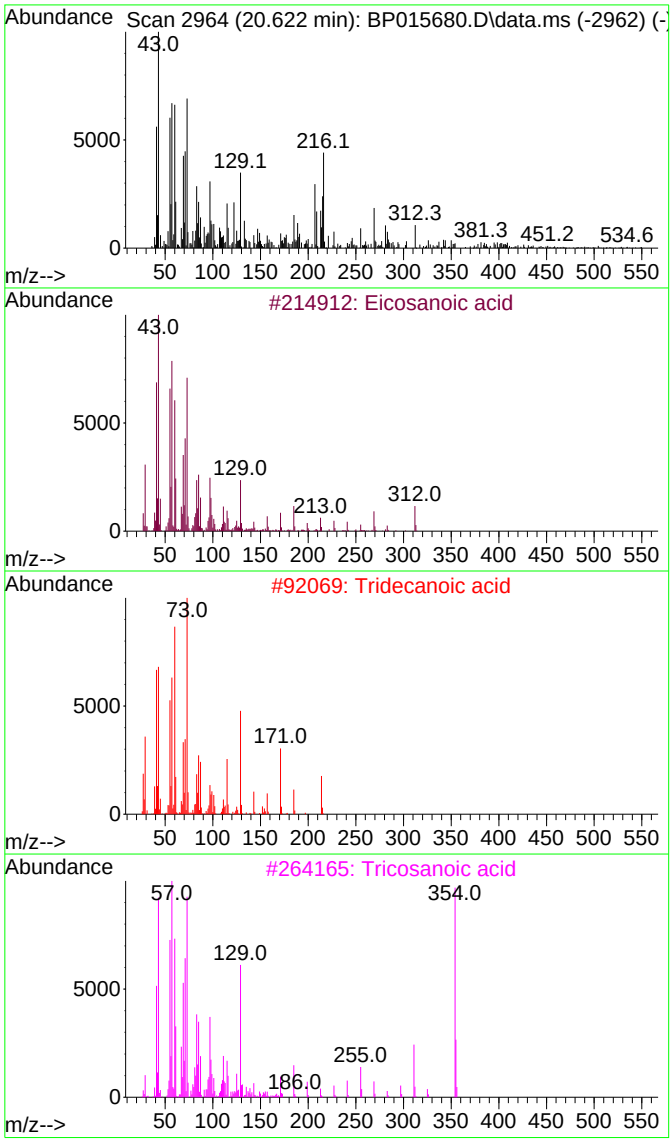
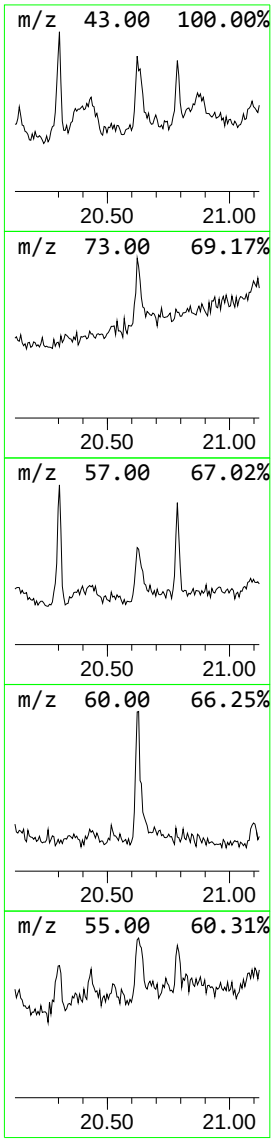
Peak Number 5 Eicosanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.622	2.00 ng/ul	171363	Chrysene-d12	21.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid	312	C20H40O2	000506-30-9	74
2	Tridecanoic acid	214	C13H26O2	000638-53-9	70
3	Tricosanoic acid	354	C23H46O2	002433-96-7	62
4	Tridecanoic acid	214	C13H26O2	000638-53-9	52
5	Octadecanoic acid	284	C18H36O2	000057-11-4	48





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

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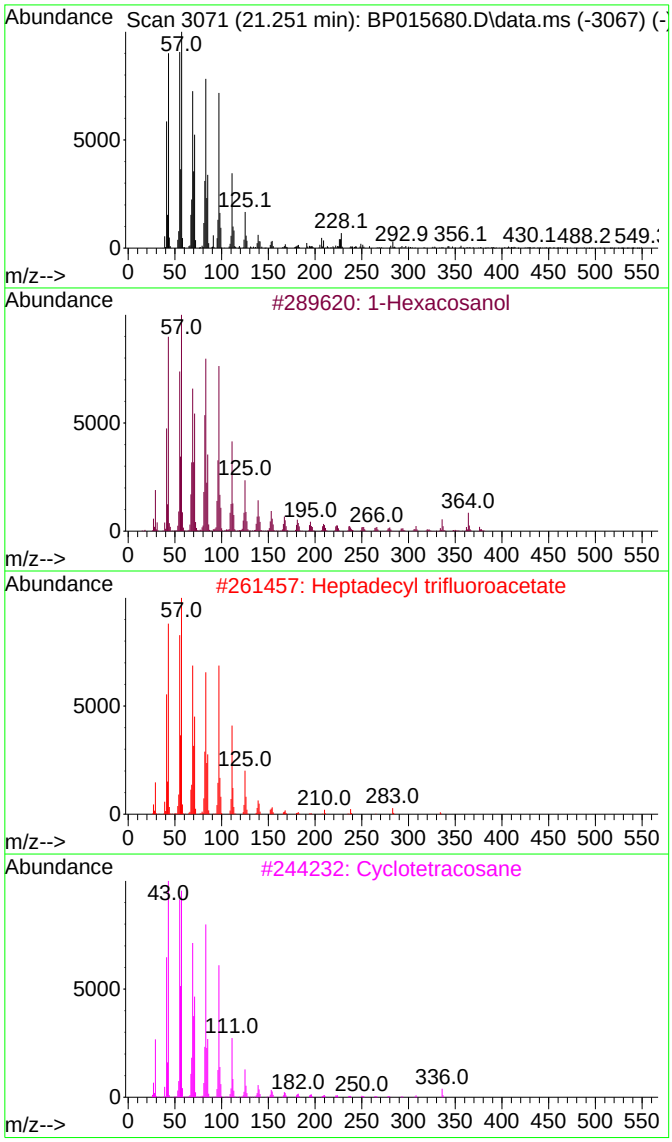
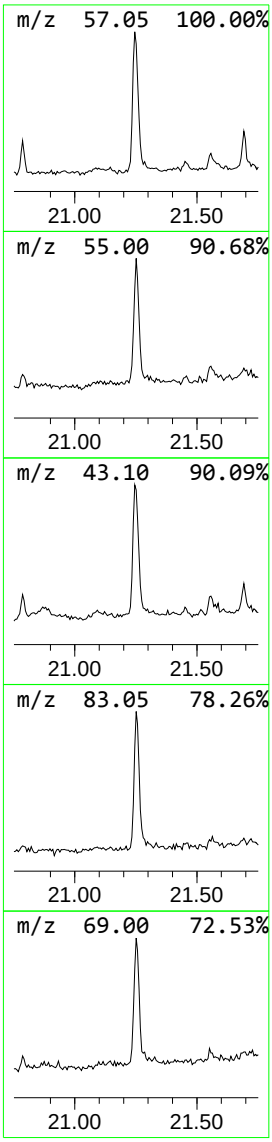
Peak Number 6 1-Hexacosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	8.03 ng/ul	687639	Chrysene-d12	21.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosanol	382	C26H54O	000506-52-5	94
2	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
3	Cyclotetracosane	336	C24H48	000297-03-0	94
4	1-Hexacosene	364	C26H52	018835-33-1	94
5	Pentacos-1-ene	350	C25H50	016980-85-1	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

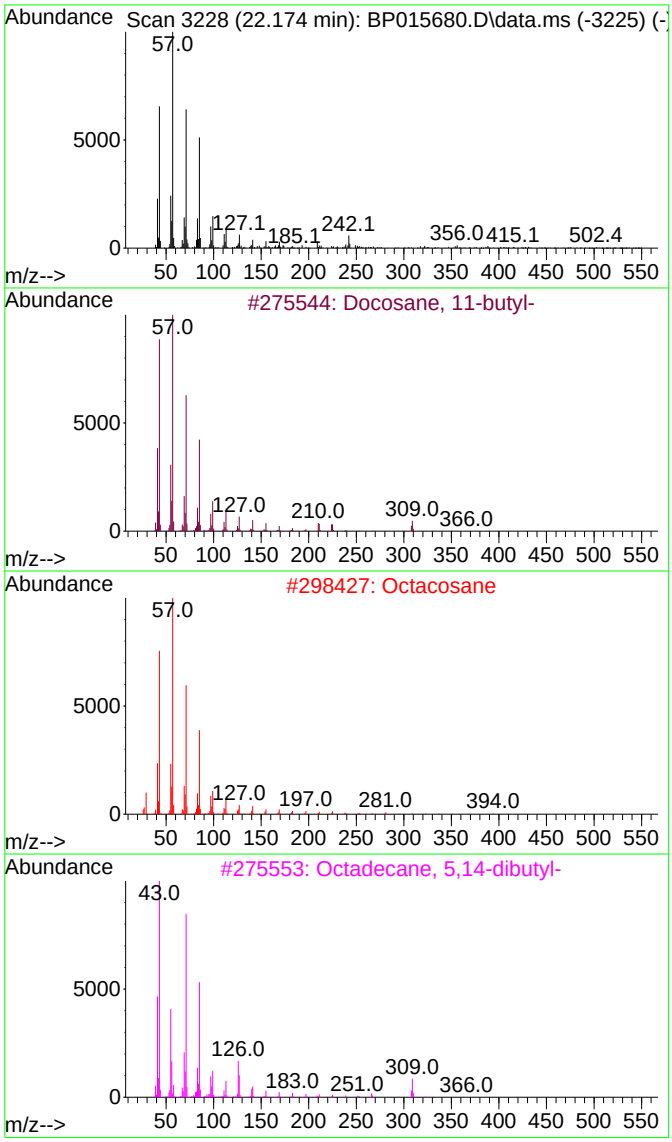
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

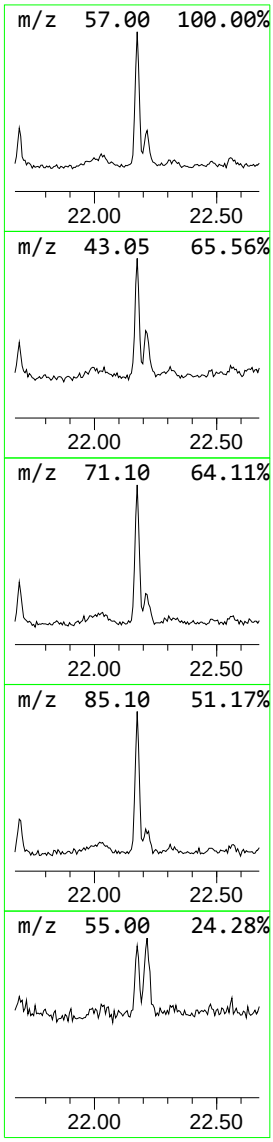
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	3.02 ng/ul	258713	Chrysene-d12	21.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2	Octacosane	394	C28H58	000630-02-4	92
3	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	90
4	Hexacosane	366	C26H54	000630-01-3	87
5	2-Methylhentriacontane	451	C32H66	001720-12-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

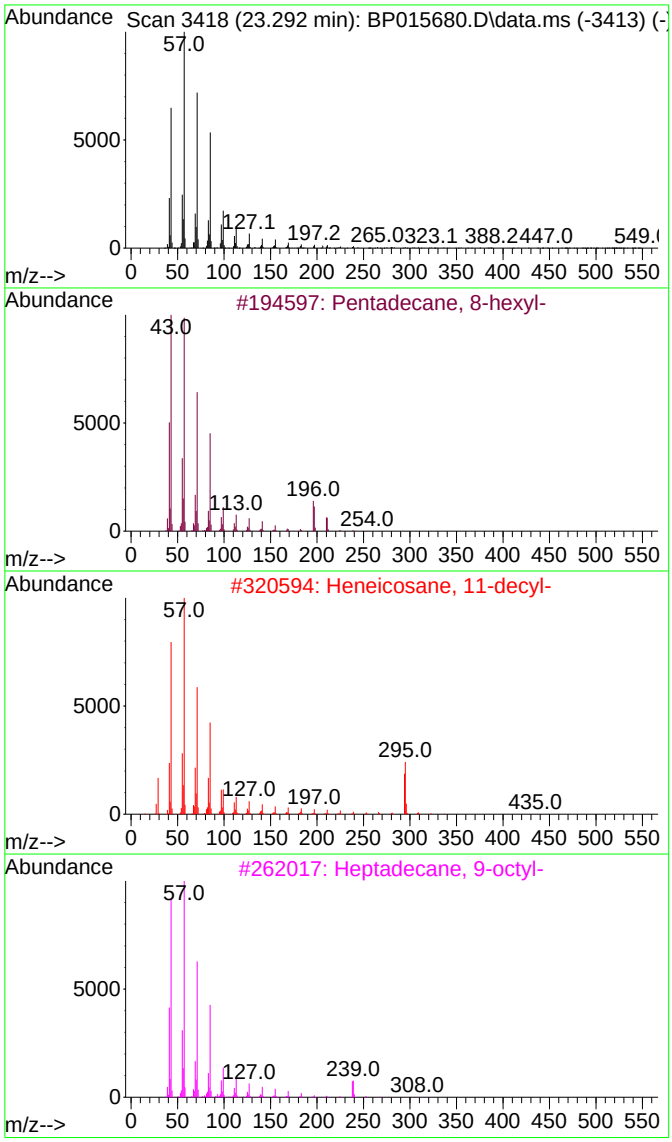
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

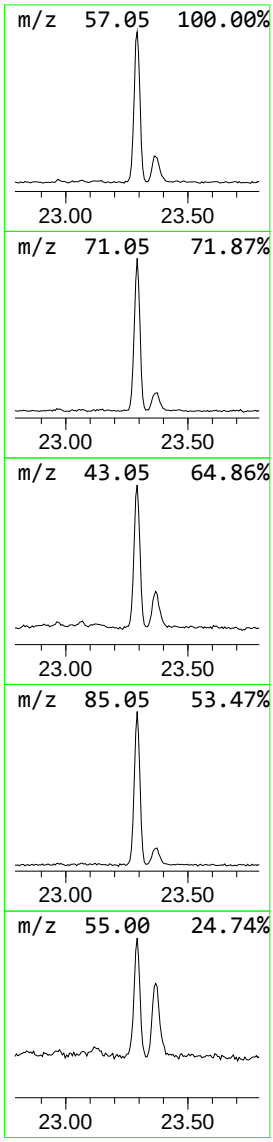
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.292	15.70 ng/ul	1400830	Perylene-d12	24.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	92
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

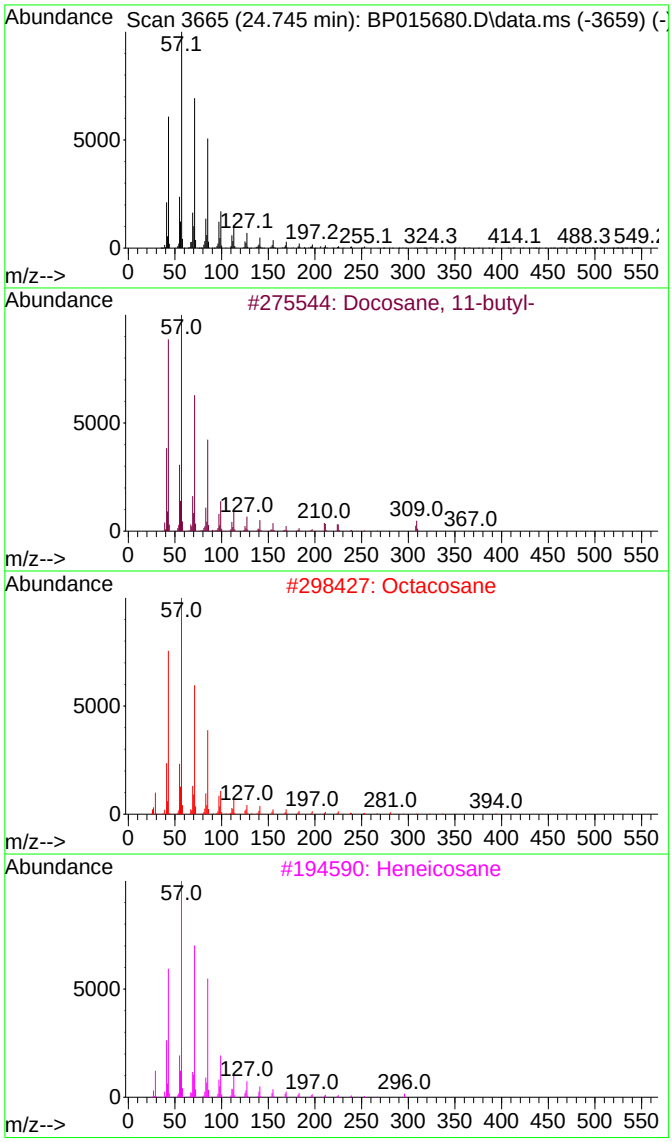
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

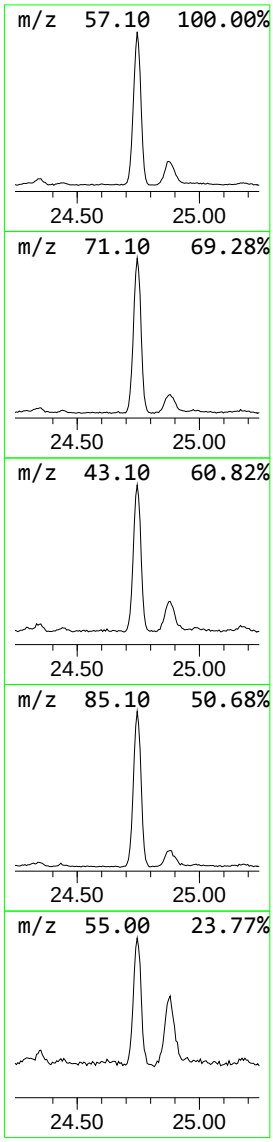
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.745	25.14 ng/ul	2243300	Perylene-d12	24.145

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Hexacosane	366	C26H54	000630-01-3	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015680.D
Acq On : 23 Jun 2023 20:01
Operator : MA/JU
Sample : 03084-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
Aceto phenone	8.922	3.5	ng/ul	166265	1	8.081	942273	20.0
Benzophenone	16.092	3.1	ng/ul	215992	3	14.728	1396080	20.0
n-Hexadecanoic ...	18.363	111.7	ng/ul	8468480	4	17.492	1517010	20.0
Octadecanoic acid	19.581	37.7	ng/ul	3228460	5	21.580	1712230	20.0
Eicosanoic acid	20.622	2.0	ng/ul	171363	5	21.580	1712230	20.0
1-Hexacosanol	21.251	8.0	ng/ul	687639	5	21.580	1712230	20.0
(DEL) Alkane: S...	22.174	3.0	ng/ul	258713	5	21.580	1712230	20.0
(DEL) Alkane: S...	23.292	15.7	ng/ul	1400830	6	24.145	1784450	20.0
(DEL) Alkane: S...	24.745	25.1	ng/ul	2243300	6	24.145	1784450	20.0