LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4661

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_G\Methods\SFAM-EPA-BG071723.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG058293.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	40	43	45	rBV3	9403	10498	0.38%	0.022%
2	3.375	45	49	57	rVB	50977	73171	2.66%	0.156%
3	3.798	117	121	128	rVB	24706	33907	1.23%	0.072%
4	4.092	167	171	176	rVB3	7964	11343	0.41%	0.024%
5	4.403	218	224	247	rBV	1476226	2233801	81.24%	4.748%
6	5.003	319	326	339	rVB	429154	640780	23.30%	1.362%
7	5.619	425	431	436	rBV2	6722	11124	0.40%	0.024%
8	5.790	456	460	466	rVB5	7701	12548	0.46%	0.027%
9	6.084	505	510	514	rBV	8856	14235	0.52%	0.030%
10	7.041	665	673	675	rBV	241092	449941	16.36%	0.956%
11	7.088	678	681	697	rVV	272307	479786	17.45%	1.020%
12	7.223	697	704	717	rVB	144485	247762	9.01%	0.527%
13	7.429	732	739	742	rBV	176260	317218	11.54%	0.674%
14	7.458	742	744	754	rVB	157564	248108	9.02%	0.527%
15	7.899	812	819	826	rBV	133503	226975	8.25%	0.482%
16	8.604	932	939	949	rBV	216890	384358	13.98%	0.817%
17	8.698	949	955	966	rVB	183531	307171	11.17%	0.653%
18	9.062	1009	1017	1020	rBV	169637	311958	11.35%	0.663%
19	9.104	1020	1024	1035	rVB	136433	246803	8.98%	0.525%
20	9.785	1132	1140	1142	rBV	139828	258332	9.40%	0.549%
21	10.191	1203	1209	1216	rBV4	4410	8558	0.31%	0.018%
22	10.326	1224	1232	1234	rBV	232037	444615	16.17%	0.945%
23	10.702	1284	1296	1300	rBV	215653	397304	14.45%	0.844%
24	10.749	1300	1304	1312	rVV	197846	362348	13.18%	0.770%
25	10.837	1312	1319	1333	rVV	193177	373650	13.59%	0.794%
26	11.031	1345	1352	1360	rVB	209387	353983	12.87%	0.752%
27	11.983	1507	1514	1528	rBV	384076	679533	24.71%	1.444%
28	12.223	1544	1555	1560	rBV2	10645	18596	0.68%	0.040%
29	12.582	1608	1616	1626	rVV	342810	575568	20.93%	1.223%
30	12.729	1633	1641	1652	rVB	232428	387665	14.10%	0.824%
31	12.970	1675	1682	1687	rVV	370396	613441	22.31%	1.304%
32	13.011	1687	1689	1702	rVV	81607	126059	4.58%	0.268%
33	13.369	1742	1750	1754	rBV	486582	779529	28.35%	1.657%
34	13.416	1754	1758	1766	rVB	245019	393972	14.33%	0.837%
35	13.904	1835	1841	1842	rBV	8125	10872	0.40%	0.023%
36	13.939	1842	1847	1851	rVV	516414	742287	27.00%	1.578%

LSC Area Percent Report

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Misc :
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Instrument :
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ClientSampleId :
A4661

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_G\Methods\SFAM-EPA-BG071723.MA.M
Title : SVOA CALIBRATION

37	13.986	1851	1855	1864	rVV	353972	516874	18.80%	1.099%
38	14.109	1870	1876	1885	rVV	265699	413893	15.05%	0.880%
39	14.233	1887	1897	1906	rVB2	505860	820053	29.82%	1.743%
40	14.538	1942	1949	1954	rBV	387076	602719	21.92%	1.281%
41	14.603	1954	1960	1972	rVB	437715	698540	25.40%	1.485%
42	14.750	1975	1985	2004	rBV3	299664	708090	25.75%	1.505%
43	14.932	2011	2016	2028	rVB	324160	518197	18.85%	1.101%
44	15.161	2049	2055	2072	rBV	295414	457524	16.64%	0.972%
45	15.326	2077	2083	2086	rVV2	5666	9204	0.33%	0.020%
46	15.373	2086	2091	2096	rVB2	8763	12810	0.47%	0.027%
47	15.531	2111	2118	2122	rBV	715799	1120660	40.76%	2.382%
48	15.572	2122	2125	2132	rVV	308444	451155	16.41%	0.959%
49	15.661	2132	2140	2151	rVV3	502750	976926	35.53%	2.076%
50	15.749	2151	2155	2164	rVB2	13783	26288	0.96%	0.056%
51	15.849	2166	2172	2174	rBV4	4907	8477	0.31%	0.018%
52	15.890	2174	2179	2188	rVV	91510	138381	5.03%	0.294%
53	16.454	2270	2275	2280	rVV	35451	51426	1.87%	0.109%
54	16.589	2291	2298	2304	rBV2	433203	641166	23.32%	1.363%
55	16.930	2350	2356	2369	rBV	472371	776747	28.25%	1.651%
56	17.024	2369	2372	2376	rVB3	7090	10322	0.38%	0.022%
57	17.112	2383	2387	2393	rVB	34200	44121	1.60%	0.094%
58	17.182	2395	2399	2406	rVB4	8393	12163	0.44%	0.026%
59	17.282	2406	2416	2420	rBV	538083	899108	32.70%	1.911%
60	17.323	2420	2423	2428	rVV	559972	769262	27.98%	1.635%
61	17.382	2428	2433	2436	rVV2	858605	1367097	49.72%	2.906%
62	17.417	2436	2439	2451	rVV	590963	860455	31.29%	1.829%
63	17.558	2459	2463	2467	rVV2	6621	10593	0.39%	0.023%
64	17.688	2477	2485	2491	rVV	459161	739755	26.90%	1.572%
65	17.758	2491	2497	2517	rVV	482770	771903	28.07%	1.641%
66	17.893	2517	2520	2522	rVV4	7931	10510	0.38%	0.022%
67	17.929	2522	2526	2533	rVV4	13003	22035	0.80%	0.047%
68	18.164	2561	2566	2573	rBV	122823	200203	7.28%	0.426%
69	18.240	2573	2579	2590	rVB	532255	757827	27.56%	1.611%
70	18.457	2613	2616	2622	rBV5	5578	9069	0.33%	0.019%
71	18.587	2634	2638	2643	rVB2	11370	14658	0.53%	0.031%
72	18.698	2653	2657	2665	rVB4	10892	18734	0.68%	0.040%
73	19.086	2720	2723	2732	rVB6	8997	13412	0.49%	0.029%
74	19.233	2741	2748	2753	rBV3	13173	20962	0.76%	0.045%
75	19.339	2756	2766	2776	rVV	1386375	2075166	75.47%	4.411%
76	19.439	2778	2783	2789	rBV4	19865	32212	1.17%	0.068%

LSC Area Percent Report

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Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : A4661

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_G\Methods\SFAM-EPA-BG071723.MA.M
Title : SVOA CALIBRATION

77	19.580	2802	2807	2814	rVB2	17019	23037	0.84%	0.049%
78	19.668	2816	2822	2825	rBV	1084475	1756095	63.87%	3.733%
79	19.697	2825	2827	2839	rVV	1037138	1346120	48.96%	2.861%
80	20.191	2906	2911	2915	rBV	21256	31489	1.15%	0.067%
81	20.578	2972	2977	2982	rBV	501820	638991	23.24%	1.358%
82	20.684	2991	2995	2998	rBV2	14977	16997	0.62%	0.036%
83	21.184	3074	3080	3105	rBV3	77641	294726	10.72%	0.626%
84	21.413	3113	3119	3125	rBV	727613	950491	34.57%	2.020%
85	21.513	3129	3136	3143	rBV2	1250643	2749641	100.00%	5.844%
86	21.577	3143	3147	3161	rVV	839191	1377027	50.08%	2.927%
87	21.706	3165	3169	3174	rVB2	25656	32337	1.18%	0.069%
88	22.294	3263	3269	3277	rBV	33773	60396	2.20%	0.128%
89	22.652	3328	3330	3337	rVB8	5318	8844	0.32%	0.019%
90	22.964	3378	3383	3391	rVB2	27341	52746	1.92%	0.112%
91	23.669	3493	3503	3511	rBV	573642	1397829	50.84%	2.971%
92	23.745	3512	3516	3521	rVB2	35251	64659	2.35%	0.137%
93	24.450	3624	3636	3642	rBV2	580268	1649520	59.99%	3.506%
94	24.515	3642	3647	3658	rVV	469489	1239065	45.06%	2.634%
95	24.662	3662	3672	3688	rVB	374874	1045261	38.01%	2.222%
96	25.749	3848	3857	3870	rVB2	35152	107709	3.92%	0.229%
97	27.053	4072	4079	4090	rVB7	26339	94625	3.44%	0.201%
98	28.246	4263	4282	4286	rBV3	375329	1630373	59.29%	3.465%
99	28.628	4340	4347	4348	rBV4	13341	26998	0.98%	0.057%
100	30.537	4665	4672	4685	rVB10	14310	56526	2.06%	0.120%

Sum of corrected areas: 47047998

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ALS Vial : 4 Sample Multiplier: 1

Instrument :
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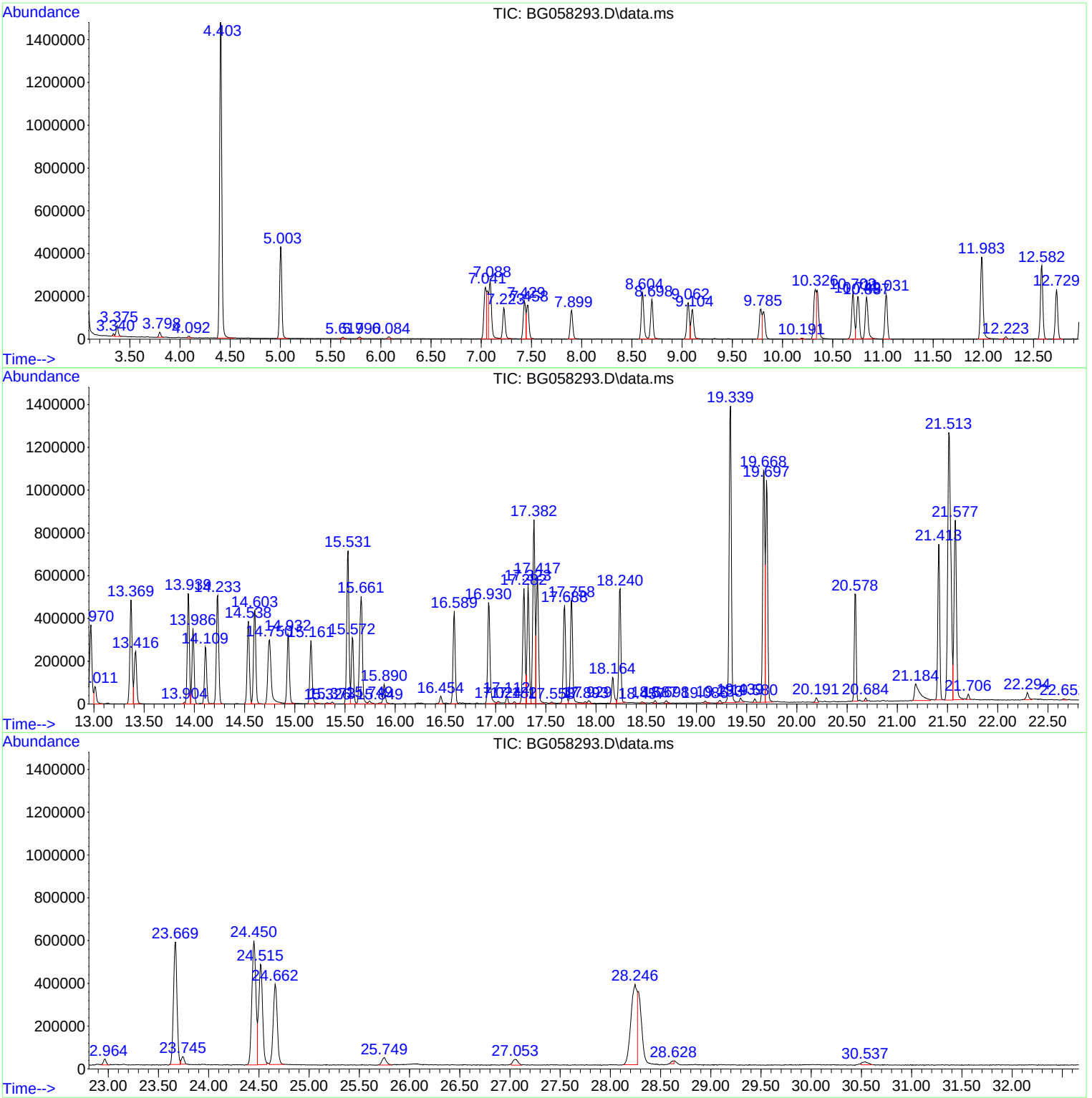
ClientSampleId :
A4661

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
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Acq On : 18 Jul 2023 8:59
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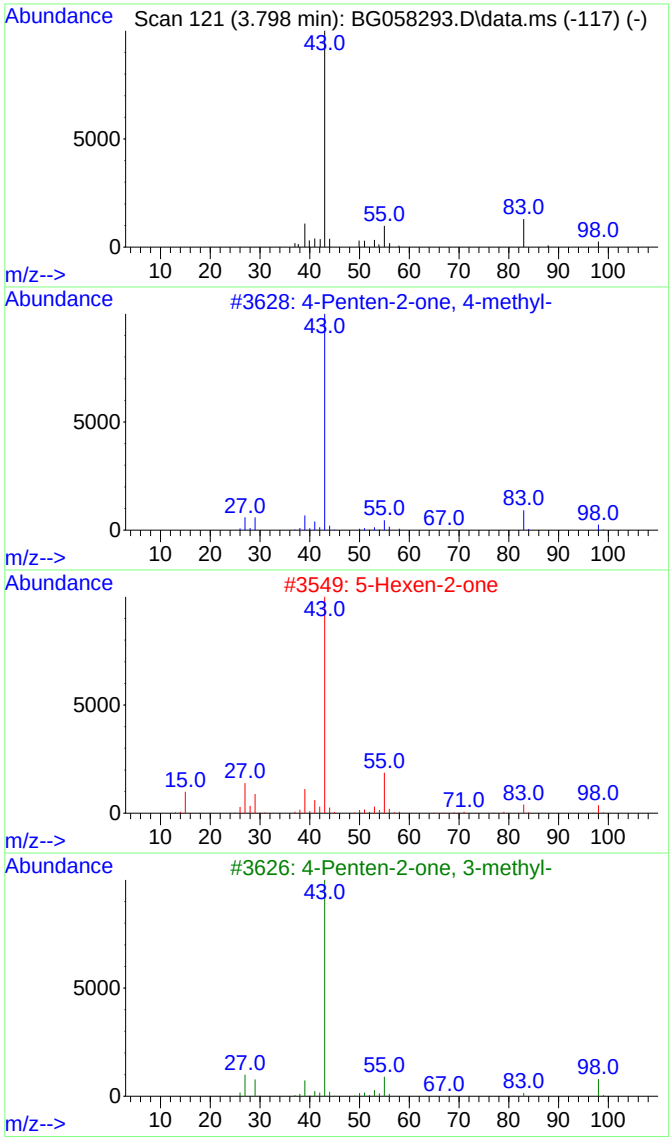
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

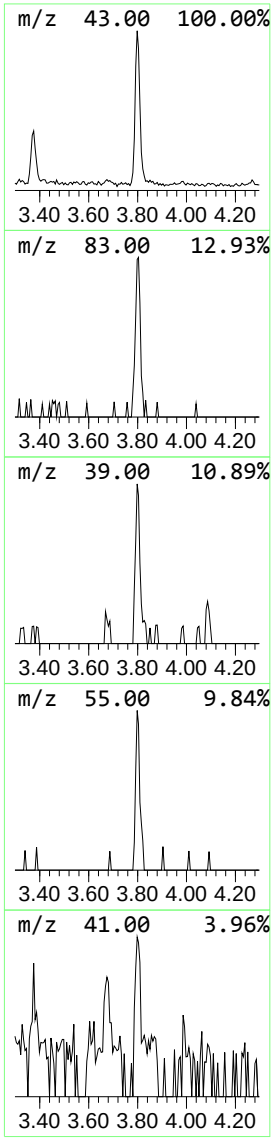
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.798	2.99 ng/ul	33907	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	64
2	5-Hexen-2-one	98	C6H10O	000109-49-9	50
3	4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	50
4	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	42
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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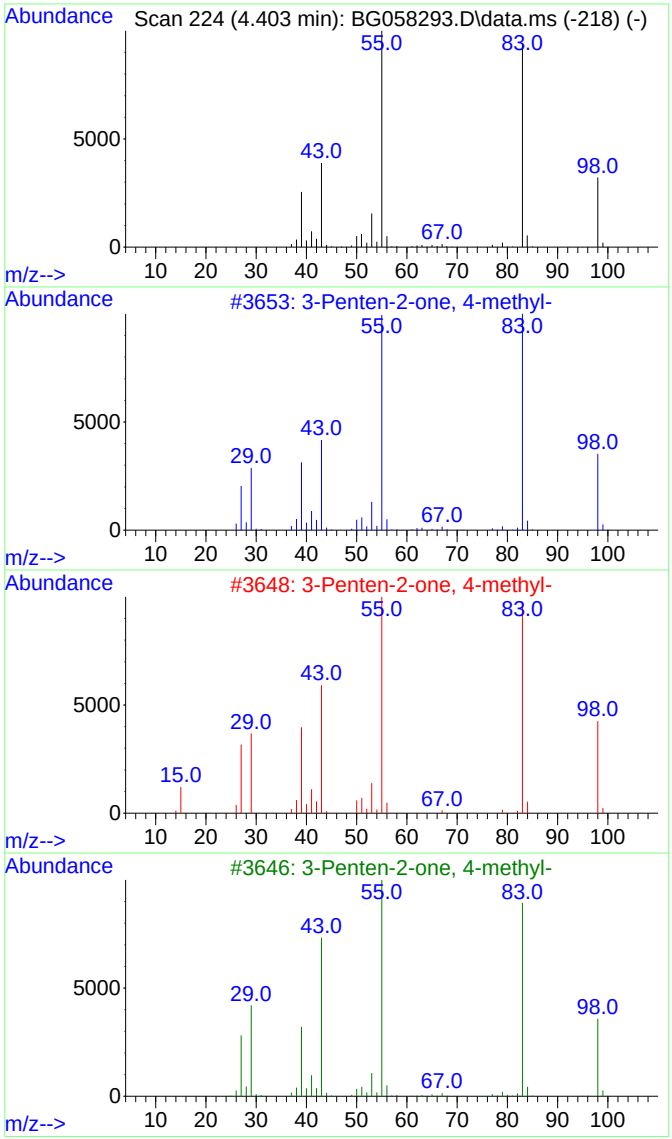
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

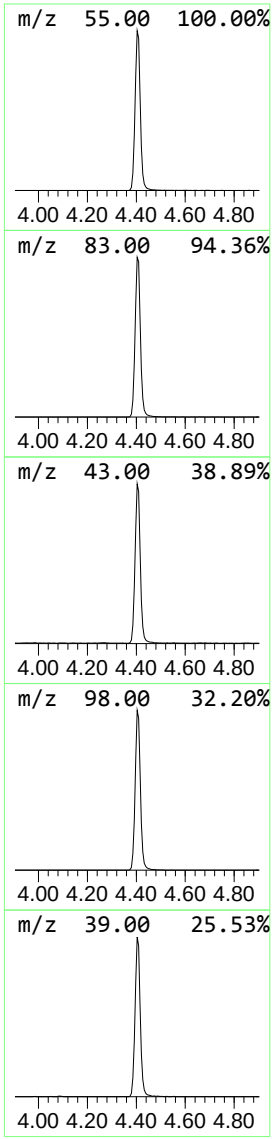
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.403	196.83 ng/ul	2233800	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
4	3-Hexen-2-one	98	C6H10O	000763-93-9	91
5	3-Hexen-2-one	98	C6H10O	000763-93-9	91





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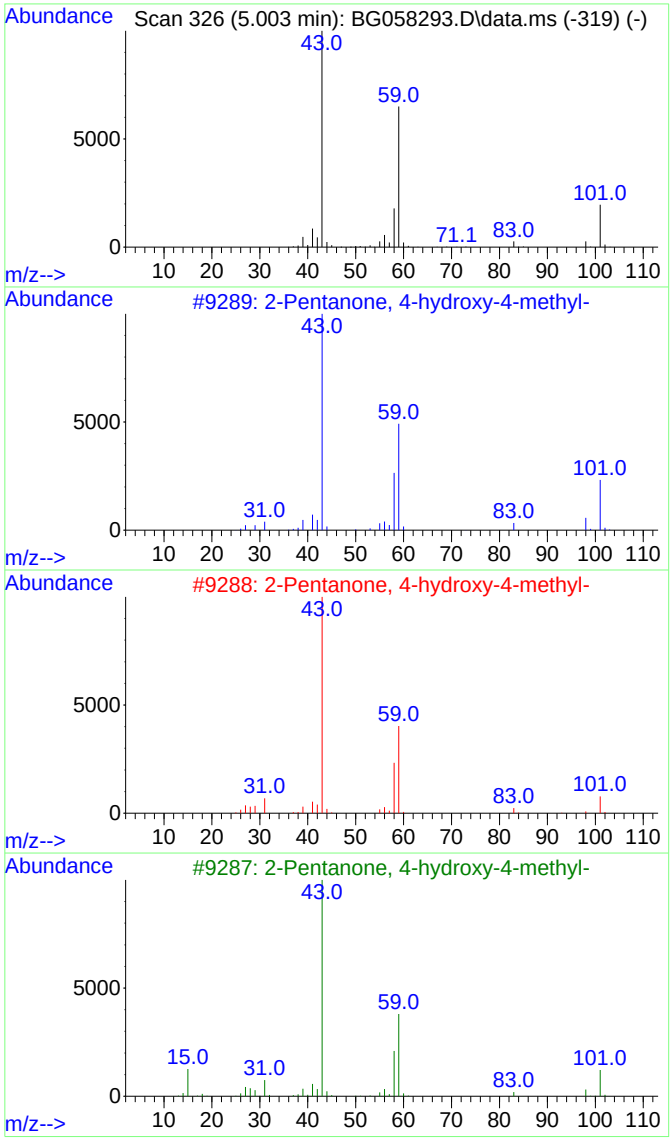
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

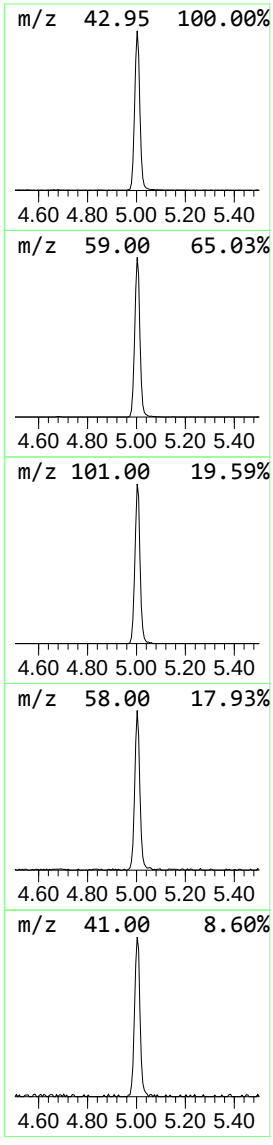
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.003	56.46 ng/ul	640780	1,4-Dichlorobenzene-d4	7.899

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36





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Instrument :
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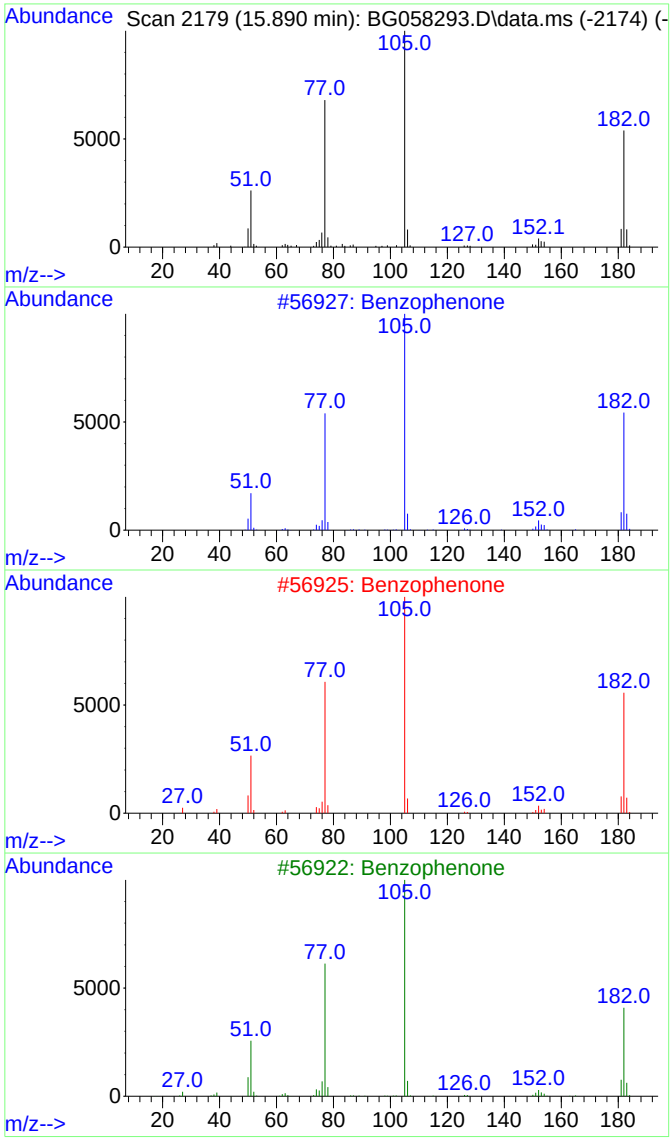
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

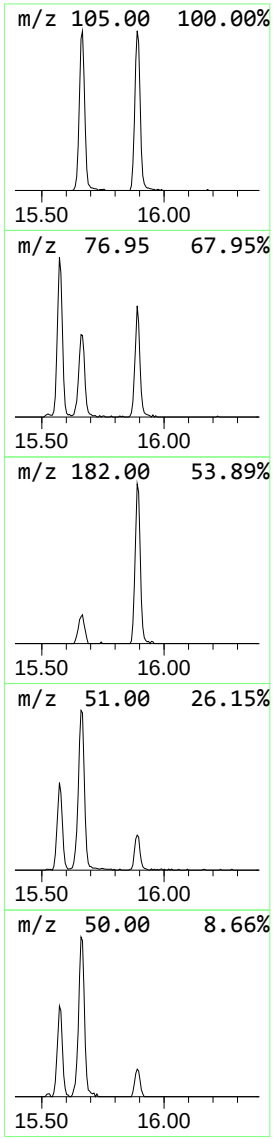
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.890	4.59 ng/ul	138381	Acenaphthene-d10	14.538

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	95
2	Benzophenone	182	C13H10O	000119-61-9	94
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	94
5	Benzophenone	182	C13H10O	000119-61-9	90





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Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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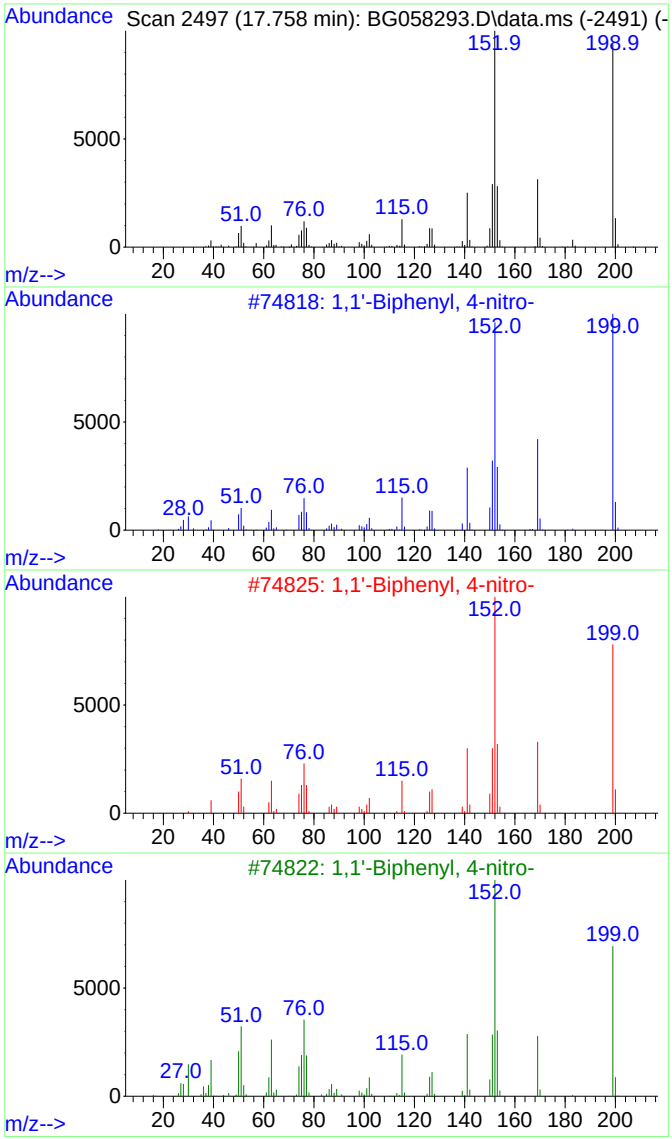
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

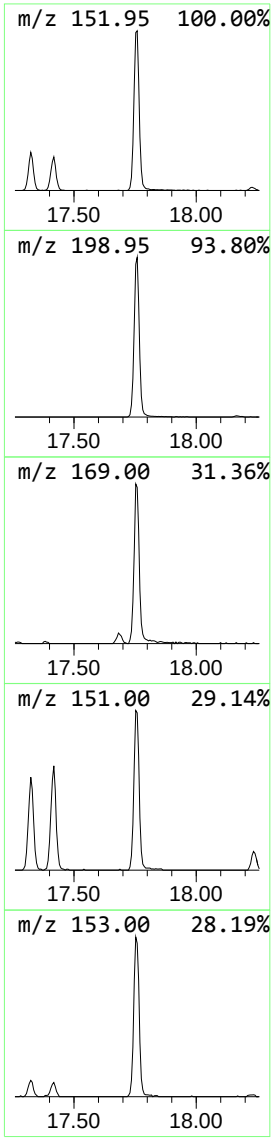
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1,1'-Biphenyl, 4-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.758	17.17 ng/ul	771903	Phenanthrene-d10	17.282

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	99
2	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	97
4	1,1'-Biphenyl, 3-nitro-	199	C12H9NO2	002113-58-8	86
5	Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	72





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Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4661

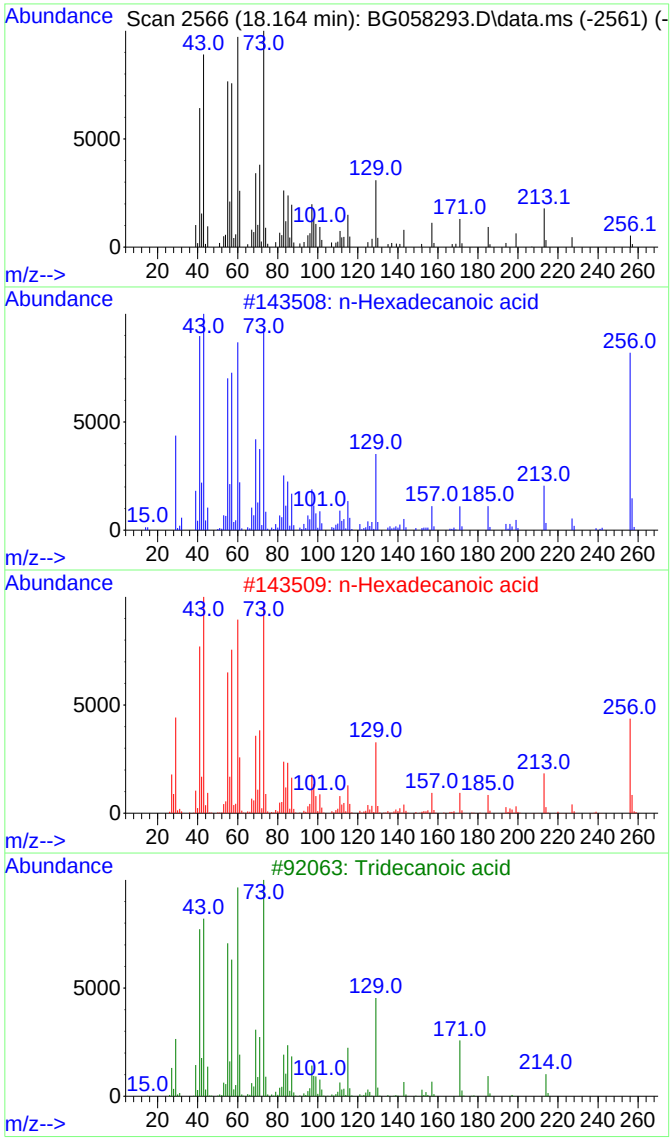
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

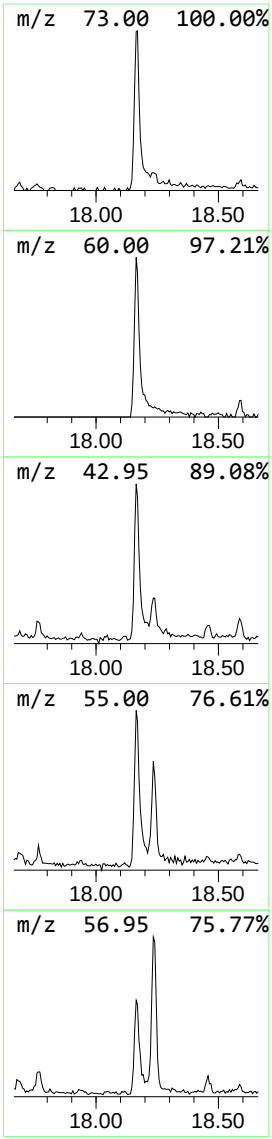
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.164	4.45 ng/ul	200203	Phenanthrene-d10	17.282

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	Tridecanoic acid	214	C13H26O2	000638-53-9	94
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4661

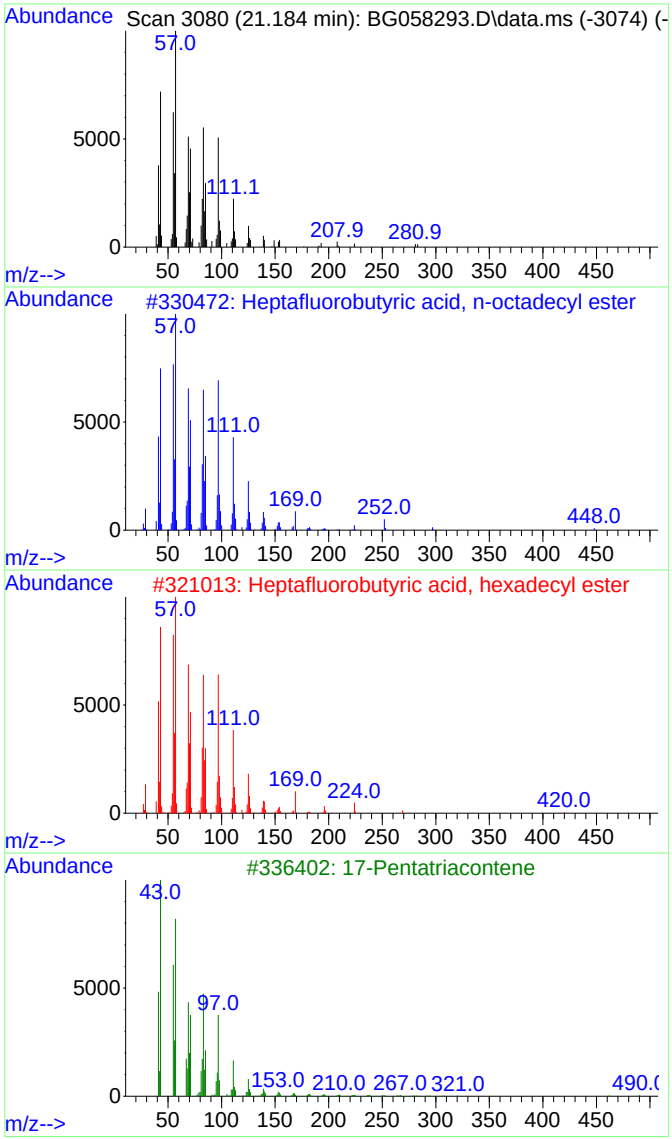
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

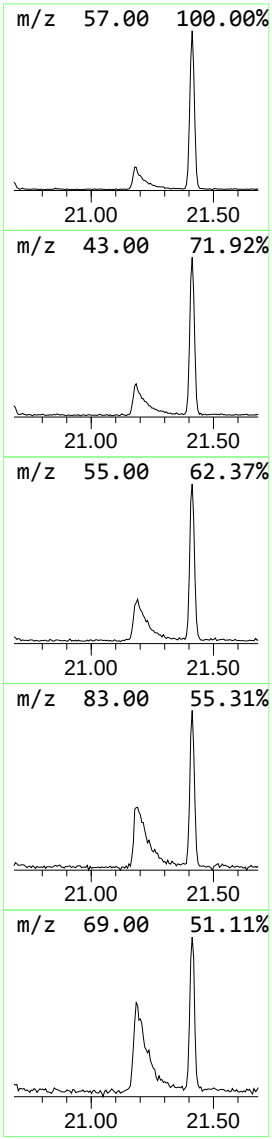
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Heptafluorobutyric acid, n-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.184	2.14 ng/ul	294726	Chrysene-d12	21.530

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
2	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
3	17-Pentatriacontene	491	C35H70	006971-40-0	91
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4661

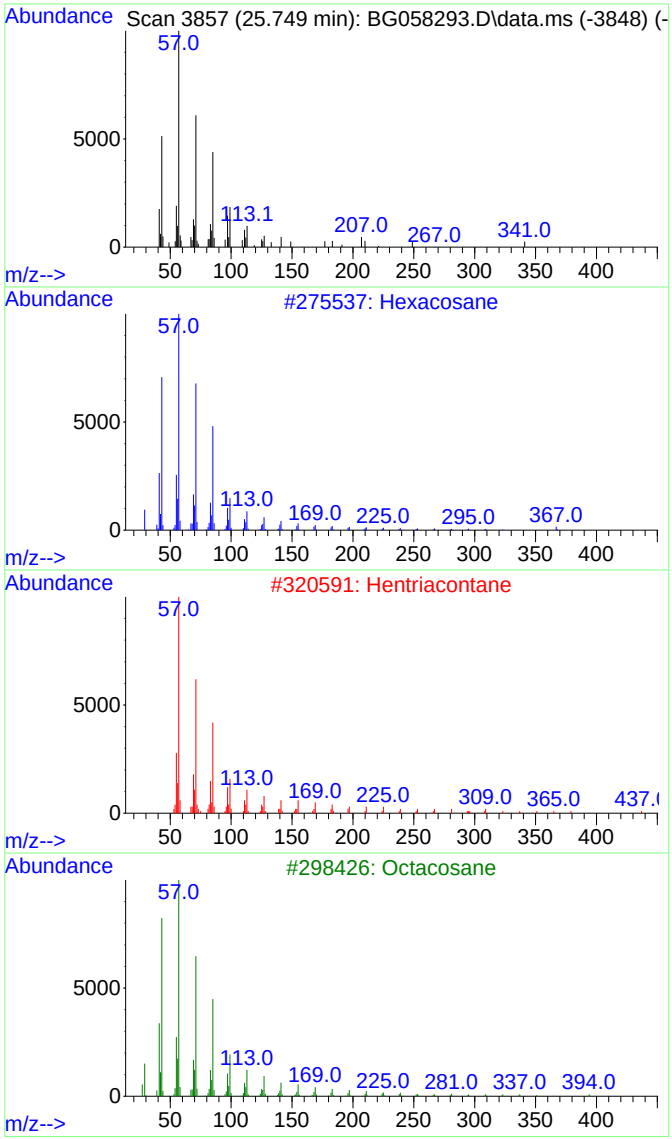
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

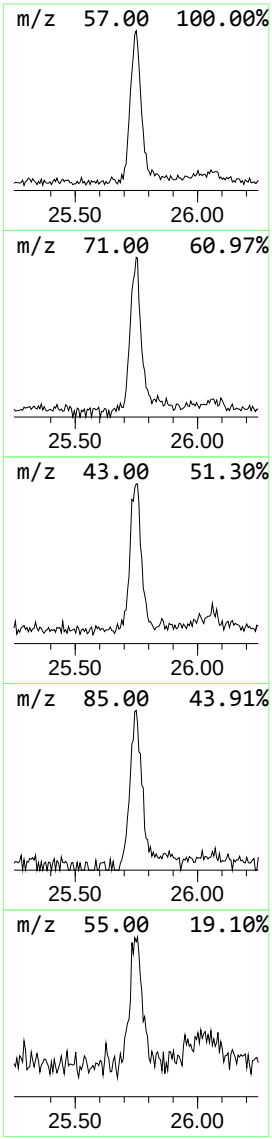
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.749	2.06 ng/ul	107709	Perylene-d12	24.662

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	Hentriacontane	437	C31H64	000630-04-6	87
3	Octacosane	394	C28H58	000630-02-4	86
4	Octacosane	394	C28H58	000630-02-4	86
5	Tetracosane	338	C24H50	000646-31-1	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058293.D
Acq On : 18 Jul 2023 8:59
Operator : CG/JU
Sample : 03458-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4661

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.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
4-Penten-2-one,...	3.798	3.0	ng/ul	33907	1	7.899	226975	20.0
3-Penten-2-one,...	4.403	196.8	ng/ul	2233800	1	7.899	226975	20.0
2-Pentanone, 4-...	5.003	56.5	ng/ul	640780	1	7.899	226975	20.0
Benzophenone	15.890	4.6	ng/ul	138381	3	14.538	602719	20.0
1,1'-Biphenyl, ...	17.758	17.2	ng/ul	771903	4	17.282	899108	20.0
n-Hexadecanoic ...	18.164	4.5	ng/ul	200203	4	17.282	899108	20.0
Heptafluorobuty...	21.184	2.1	ng/ul	294726	5	21.530	2749640	20.0
(DEL) Alkane: S...	25.749	2.1	ng/ul	107709	6	24.662	1045260	20.0