LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

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-BP062623.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP015984.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.258	9	12	20	rVB3	30409	52790	0.78%	0.047%
2	3.370	28	31	36	rVB	62044	77996	1.15%	0.069%
3	3.817	103	107	112	rBV	31705	40130	0.59%	0.036%
4	3.858	112	114	119	rVV	23297	36999	0.55%	0.033%
5	3.917	120	124	133	rVV	169593	236520	3.50%	0.209%
6	3.993	133	137	139	rVV2	24986	35271	0.52%	0.031%
7	4.034	139	144	149	rVV	127564	197987	2.93%	0.175%
8	4.134	158	161	168	rVB	36029	52717	0.78%	0.047%
9	4.487	216	221	226	rBV	97569	154459	2.28%	0.137%
10	5.105	318	326	334	rBV	290836	494340	7.31%	0.438%
11	5.999	473	478	485	rBV8	17954	36376	0.54%	0.032%
12	7.234	677	688	704	rBV	631596	1683630	24.90%	1.491%
13	7.369	705	711	739	rVB	734396	1415418	20.94%	1.254%
14	7.575	739	746	759	rBV	895446	1740450	25.74%	1.541%
15	8.040	818	825	840	rBV	1876309	3476175	51.42%	3.079%
16	8.593	915	919	928	rVB5	18692	42702	0.63%	0.038%
17	8.793	945	953	966	rBV	609090	1715032	25.37%	1.519%
18	8.899	966	971	985	rVB	208371	457719	6.77%	0.405%
19	9.234	1021	1028	1042	rBV	815505	1701073	25.16%	1.507%
20	9.963	1141	1152	1172	rBV	520731	1455078	21.52%	1.289%
21	10.257	1194	1202	1209	rBV6	31153	105206	1.56%	0.093%
22	10.357	1215	1219	1228	rVB2	15866	33300	0.49%	0.029%
23	10.540	1243	1250	1280	rBV	611311	1996126	29.53%	1.768%
24	10.869	1298	1306	1323	rVV	2568336	5198740	76.90%	4.604%
25	11.063	1332	1339	1359	rBV2	240907	754819	11.16%	0.669%
26	11.763	1455	1458	1468	rVV2	14819	39612	0.59%	0.035%
27	11.852	1468	1473	1483	rVB	31996	58294	0.86%	0.052%
28	12.257	1537	1542	1549	rBV3	24763	49633	0.73%	0.044%
29	12.551	1587	1592	1602	rVV	28759	62018	0.92%	0.055%
30	12.763	1621	1628	1634	rBV	28705	58755	0.87%	0.052%
31	13.193	1696	1701	1705	rBV3	29587	44614	0.66%	0.040%
32	13.487	1745	1751	1757	rBV	39740	65044	0.96%	0.058%
33	13.569	1757	1765	1770	rBV2	57014	122795	1.82%	0.109%
34	13.669	1777	1782	1786	rVB	41965	58933	0.87%	0.052%
35	13.881	1811	1818	1822	rBV6	18334	40256	0.60%	0.036%
36	14.010	1833	1840	1846	rBV4	34264	85929	1.27%	0.076%

LSC Area Percent Report

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Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

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-BP062623.MA.M
Title : SVOA CALIBRATION

37	14.104	1847	1856	1867	rVV	1736111	3024549	44.74%	2.679%
38	14.387	1898	1904	1918	rBV2	2258791	3657709	54.10%	3.240%
39	14.551	1928	1932	1940	rVB	63121	94953	1.40%	0.084%
40	14.693	1949	1956	1963	rBV	4197446	6371394	94.24%	5.643%
41	14.757	1963	1967	1977	rVB	116163	211159	3.12%	0.187%
42	15.004	2003	2009	2022	rBV3	190904	723280	10.70%	0.641%
43	15.104	2023	2026	2034	rVB3	73870	146472	2.17%	0.130%
44	15.316	2057	2062	2066	rVB6	18759	33102	0.49%	0.029%
45	15.457	2082	2086	2090	rBV5	23252	39622	0.59%	0.035%
46	15.504	2091	2094	2100	rVB2	72894	97214	1.44%	0.086%
47	15.692	2119	2126	2133	rBV	2550595	4157992	61.50%	3.683%
48	15.869	2150	2156	2175	rBV2	308210	865210	12.80%	0.766%
49	16.057	2182	2188	2199	rBV	877730	1398833	20.69%	1.239%
50	16.198	2209	2212	2215	rBV	22554	31765	0.47%	0.028%
51	16.387	2237	2244	2249	rVV3	106195	243795	3.61%	0.216%
52	16.451	2252	2255	2261	rVB7	23301	40798	0.60%	0.036%
53	16.539	2266	2270	2277	rBV5	43083	83368	1.23%	0.074%
54	16.616	2277	2283	2291	rVB2	127475	244076	3.61%	0.216%
55	16.763	2305	2308	2313	rVB4	29872	42283	0.63%	0.037%
56	16.828	2317	2319	2326	rVB5	32076	53301	0.79%	0.047%
57	16.981	2342	2345	2348	rBV3	25719	31838	0.47%	0.028%
58	17.016	2349	2351	2359	rVB5	27658	43310	0.64%	0.038%
59	17.139	2367	2372	2374	rBV2	67738	109153	1.61%	0.097%
60	17.163	2374	2376	2380	rVB3	68028	85554	1.27%	0.076%
61	17.275	2391	2395	2399	rVB	60936	81279	1.20%	0.072%
62	17.322	2399	2403	2411	rBV2	127581	191133	2.83%	0.169%
63	17.392	2411	2415	2419	rBV	87476	114635	1.70%	0.102%
64	17.451	2419	2425	2430	rBV	4311298	6624061	97.98%	5.867%
65	17.498	2430	2433	2438	rVV	1378908	1969632	29.13%	1.744%
66	17.557	2438	2443	2462	rVB2	2193165	4009516	59.31%	3.551%
67	17.881	2493	2498	2501	rBV2	188491	333389	4.93%	0.295%
68	18.057	2524	2528	2530	rBV4	45453	70477	1.04%	0.062%
69	18.081	2530	2532	2538	rVB3	70227	88763	1.31%	0.079%
70	18.198	2548	2552	2556	rVB5	78901	113648	1.68%	0.101%
71	18.251	2557	2561	2566	rBV	394582	555169	8.21%	0.492%
72	18.310	2566	2571	2578	rVV2	2769259	3815538	56.44%	3.379%
73	18.369	2578	2581	2592	rVB2	300875	605870	8.96%	0.537%
74	18.510	2600	2605	2608	rBV2	222328	352525	5.21%	0.312%
75	18.569	2613	2615	2628	rVB3	130901	296772	4.39%	0.263%
76	18.798	2651	2654	2658	rBV	104570	126743	1.87%	0.112%

LSC Area Percent Report

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Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

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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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Title : SVOA CALIBRATION

77	18.869	2662	2666	2672	rVB4	126946	228834	3.38%	0.203%
78	19.181	2716	2719	2725	rBV2	163608	315506	4.67%	0.279%
79	19.392	2753	2755	2757	rBV	101291	97881	1.45%	0.087%
80	19.422	2757	2760	2761	rVV2	176362	215740	3.19%	0.191%
81	19.457	2761	2766	2773	rVB	2886965	4084371	60.41%	3.617%
82	19.533	2775	2779	2790	rBV2	681499	979776	14.49%	0.868%
83	19.733	2811	2813	2816	rVB	119662	108752	1.61%	0.096%
84	19.780	2816	2821	2824	rBV	3179144	4527677	66.97%	4.010%
85	19.810	2824	2826	2834	rVB	2212243	2685869	39.73%	2.379%
86	20.257	2898	2902	2906	rBV3	301326	509641	7.54%	0.451%
87	20.733	2980	2983	2987	rBV	381758	445741	6.59%	0.395%
88	21.033	3031	3034	3039	rBV	344414	472914	6.99%	0.419%
89	21.192	3057	3061	3063	rBV	741789	899425	13.30%	0.797%
90	21.227	3063	3067	3071	rVB2	879360	1271697	18.81%	1.126%
91	21.404	3094	3097	3100	rBV	283523	343710	5.08%	0.304%
92	21.545	3114	3121	3125	rBV2	3790321	6760766	100.00%	5.988%
93	21.580	3125	3127	3132	rVB	1048536	1282162	18.96%	1.136%
94	22.110	3215	3217	3221	rBV	446562	483812	7.16%	0.429%
95	23.216	3400	3405	3413	rVB	1982971	3359858	49.70%	2.976%
96	23.316	3415	3422	3436	rVB2	1444918	5249733	77.65%	4.650%
97	23.863	3510	3515	3519	rBV2	824236	1749824	25.88%	1.550%
98	23.916	3520	3524	3530	rVB2	1259248	2357029	34.86%	2.088%
99	24.086	3547	3553	3560	rBV	2275664	4976458	73.61%	4.408%
100	24.639	3641	3647	3655	rVB	2041978	4442353	65.71%	3.935%

Sum of corrected areas: 112906345

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
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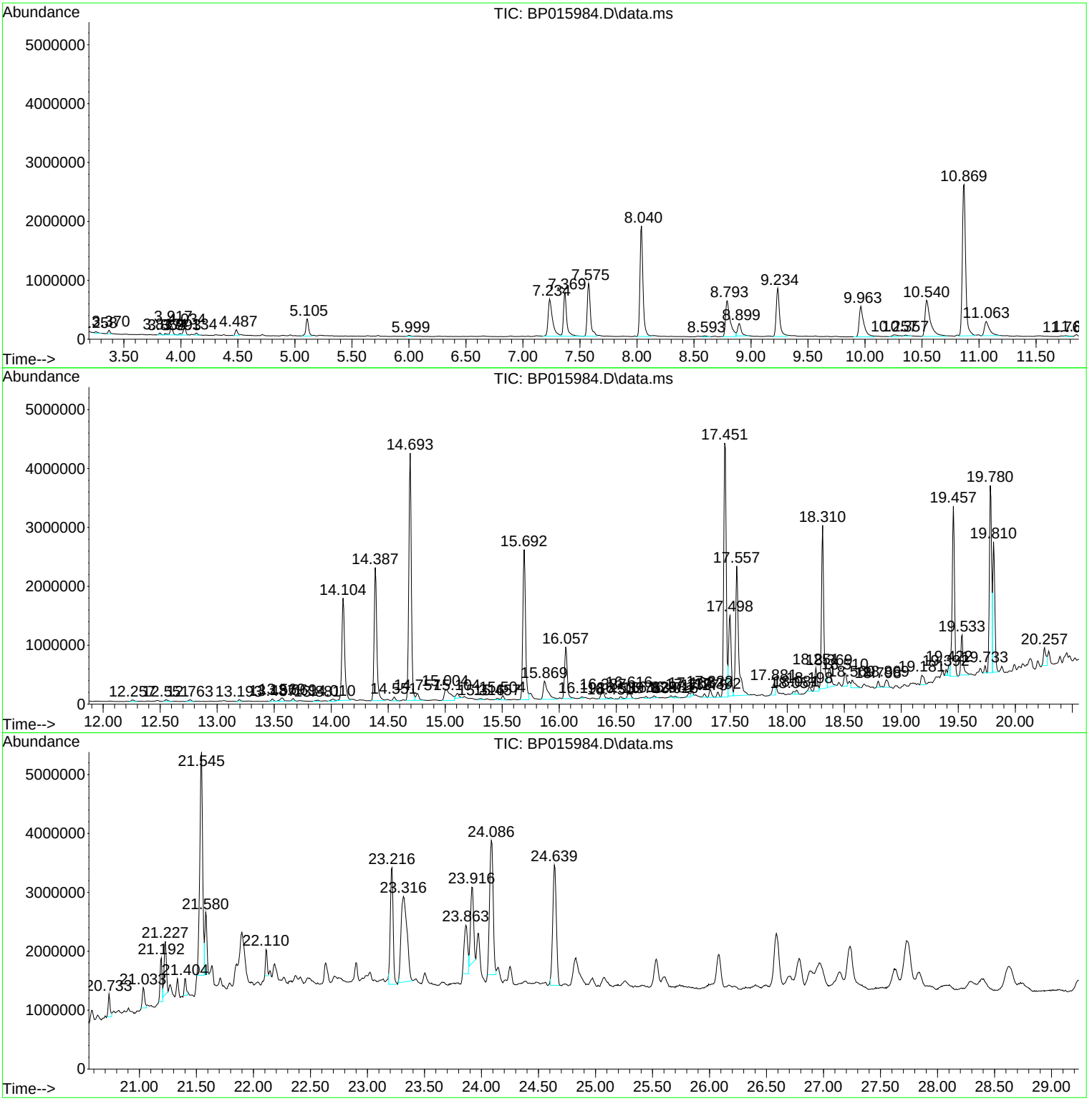
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
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ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

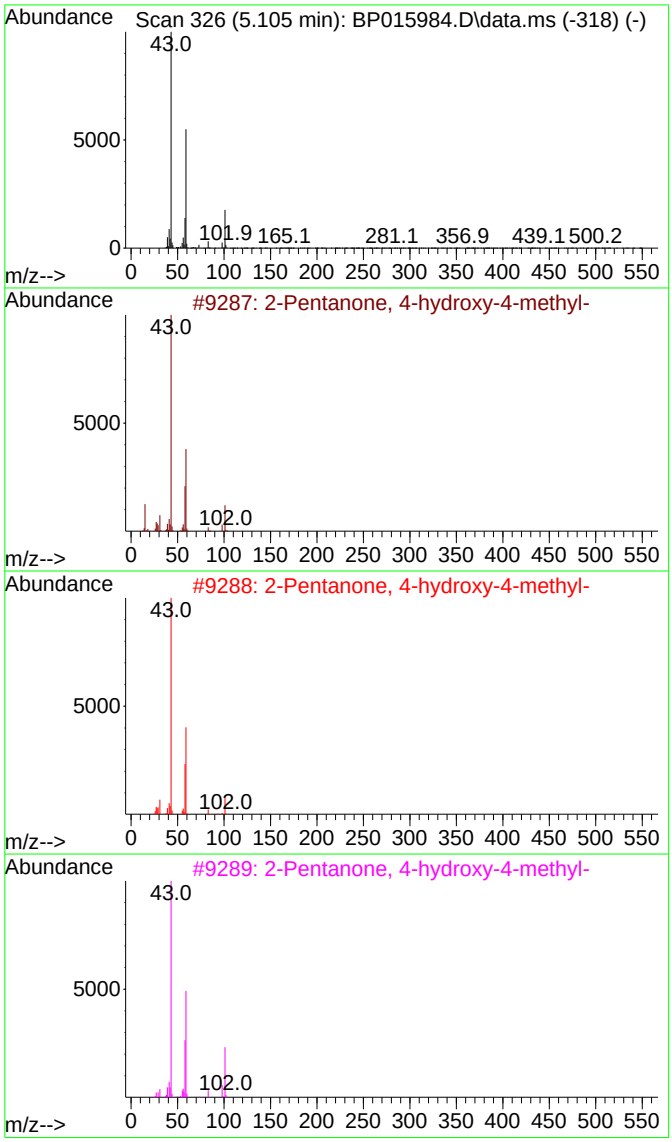
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

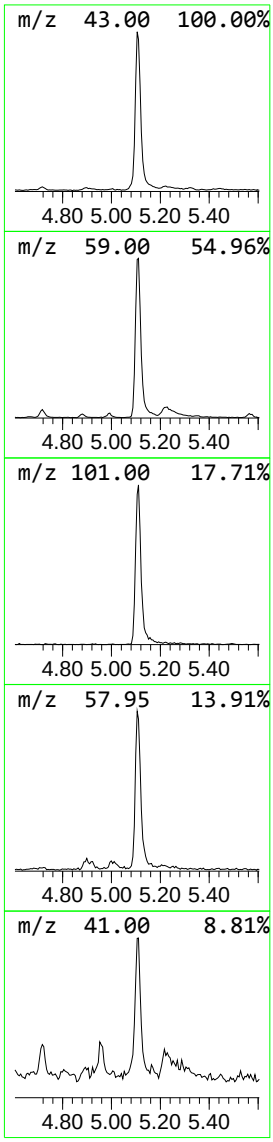
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.105	2.84 ng/ul	494340	1,4-Dichlorobenzene-d4	8.040

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
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Operator : MA/JU
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ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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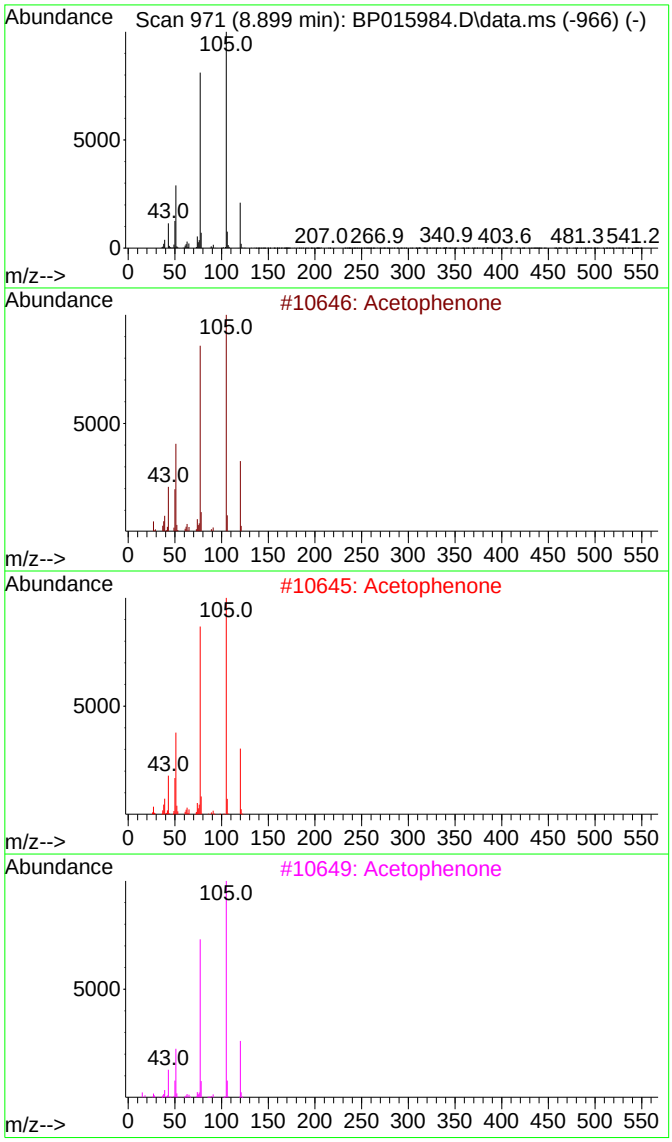
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

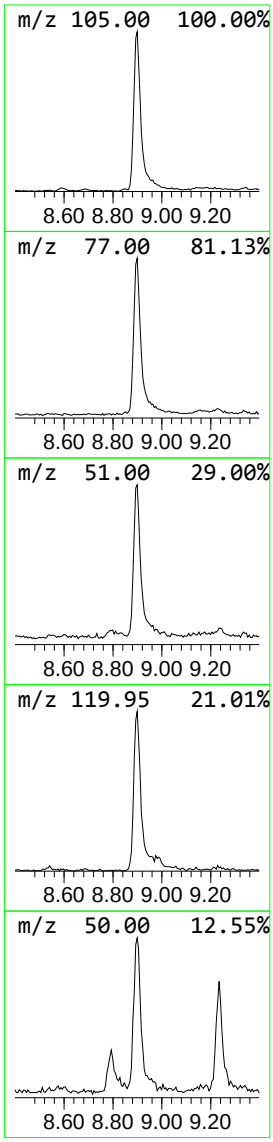
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Aceto phenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.899	2.63 ng/ul	457719	1,4-Dichlorobenzene-d4	8.040

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetophenone	120	C8H8O	000098-86-2	93
2	Acetophenone	120	C8H8O	000098-86-2	91
3	Acetophenone	120	C8H8O	000098-86-2	91
4	Acetophenone	120	C8H8O	000098-86-2	91
5	Acetophenone	120	C8H8O	000098-86-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
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ALS Vial : 30 Sample Multiplier: 1

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BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

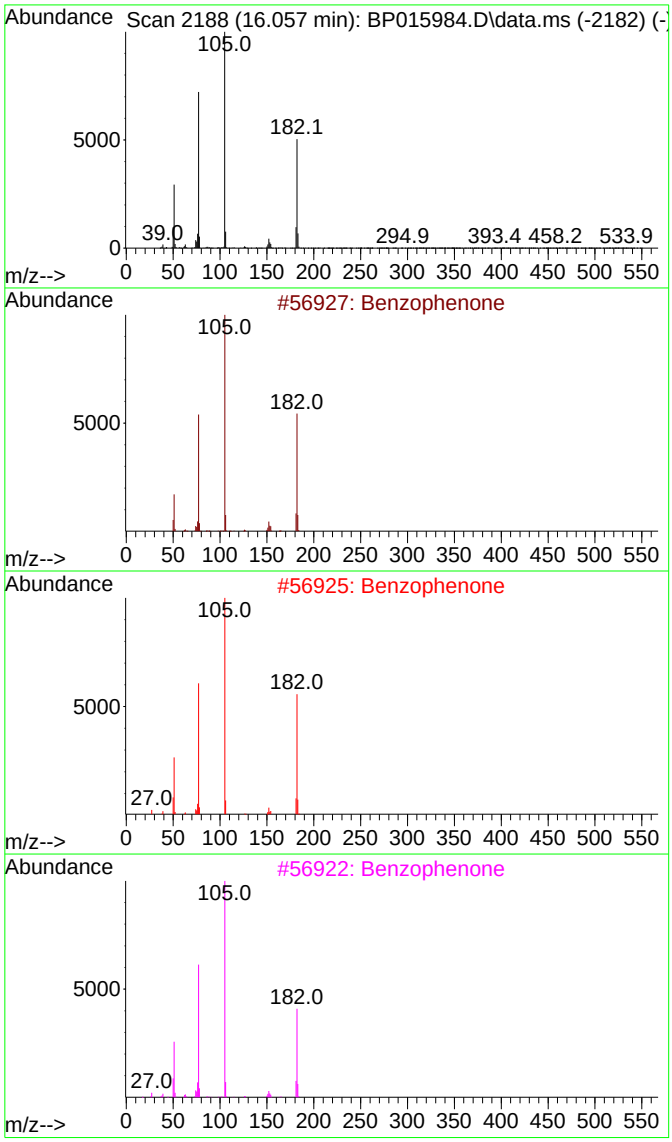
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

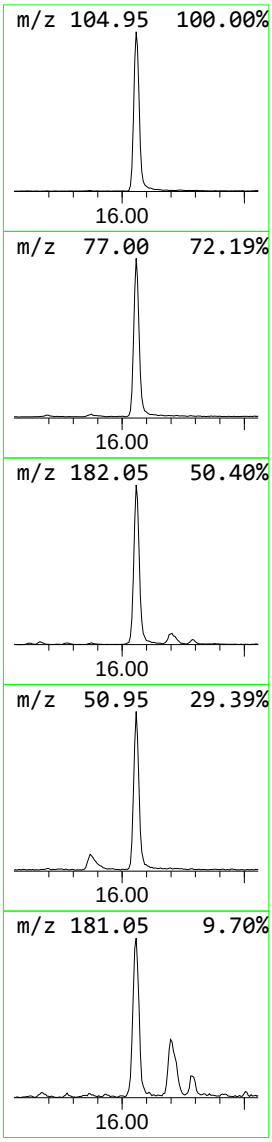
****

Peak Number 3 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.057	4.39 ng/ul	1398830	Acenaphthene-d10	14.693

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	96
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	91





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Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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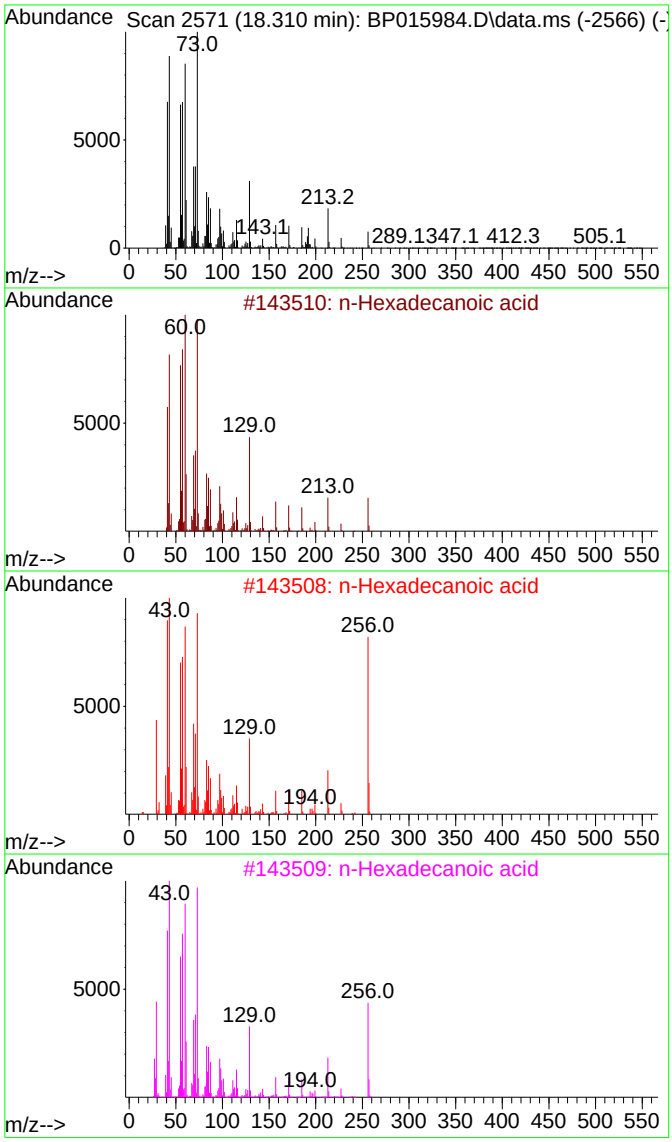
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

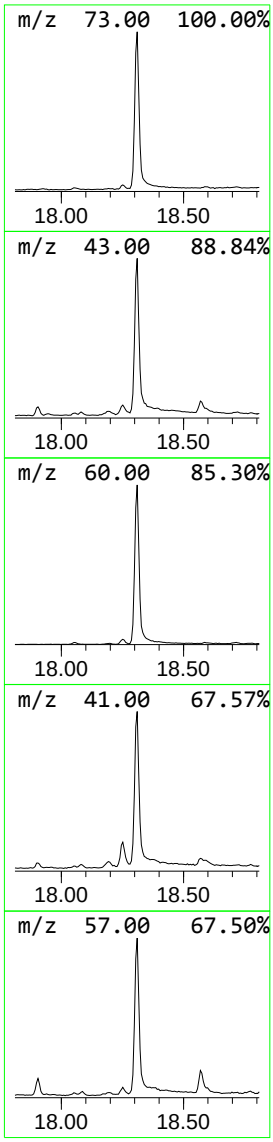
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.310	11.52 ng/ul	3815540	Phenanthrene-d10	17.451

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





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

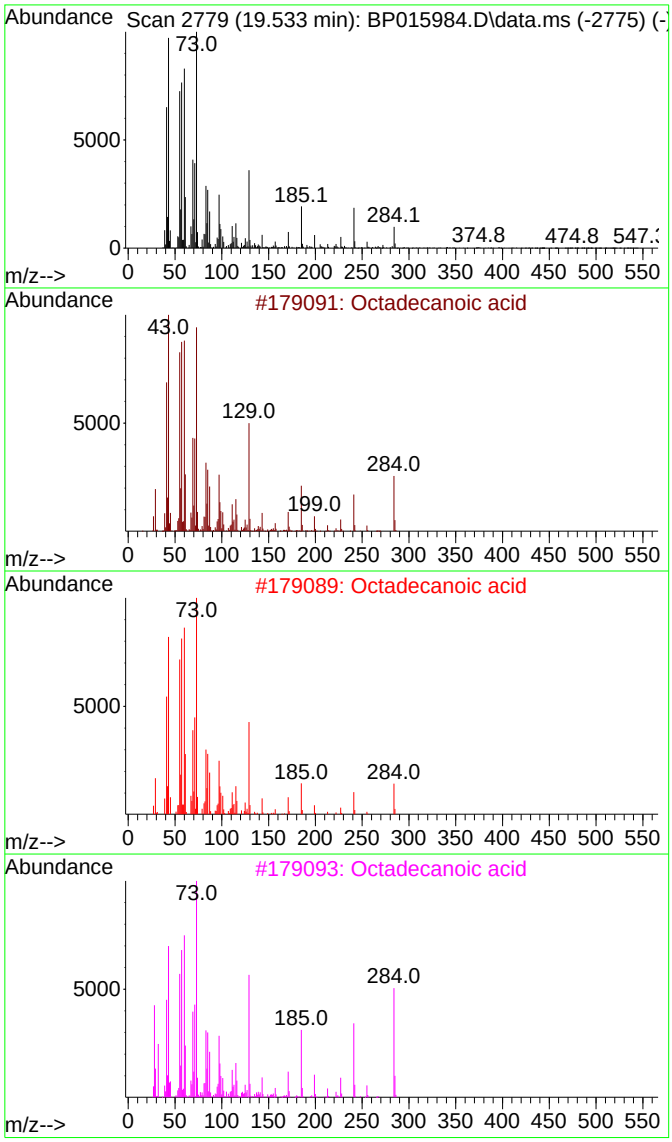
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

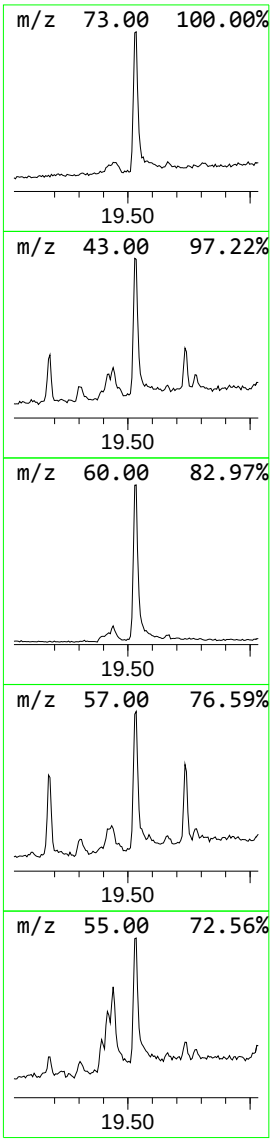
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.533	2.90 ng/ul	979776	Chrysene-d12	21.545

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	Octadecanoic acid	284	C18H36O2	000057-11-4	95





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Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

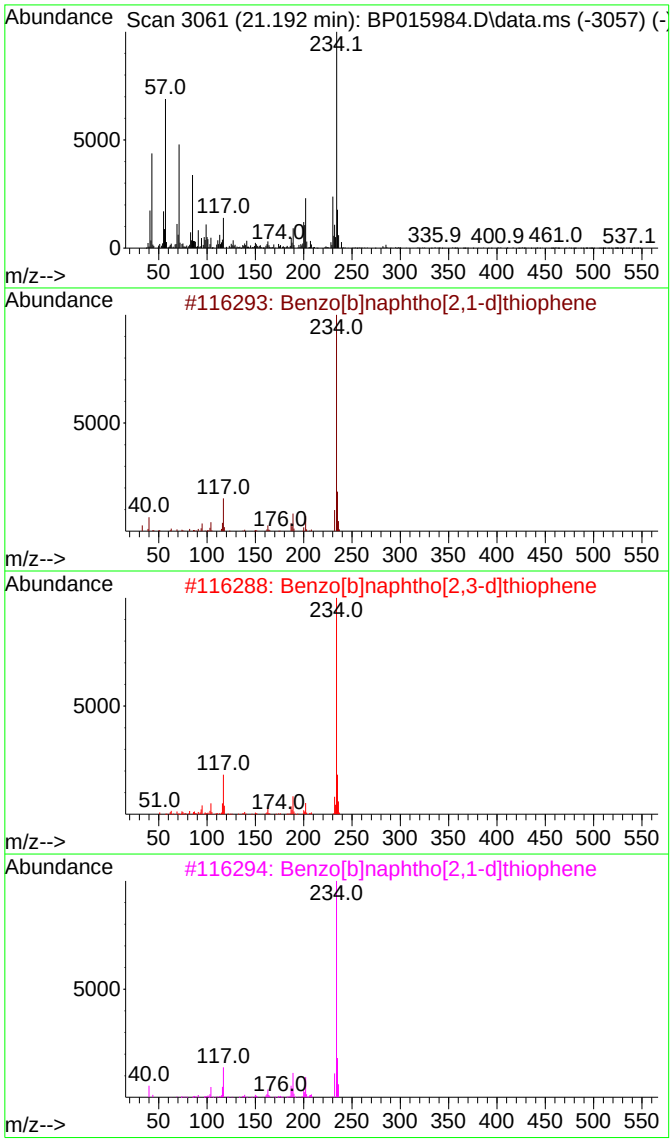
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

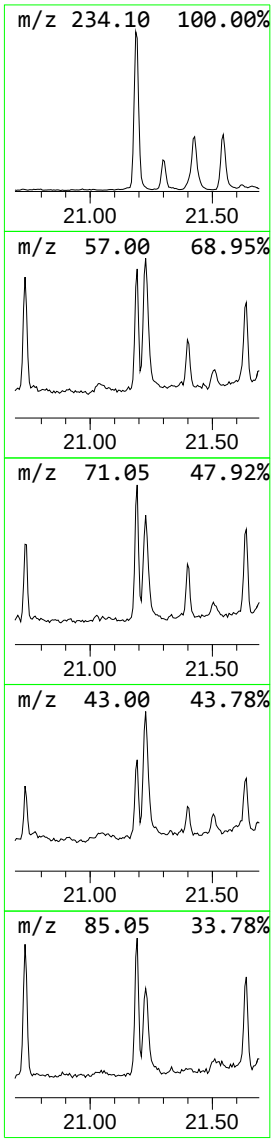
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	2.66 ng/ul	899425	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	89
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	89
3	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	89
4	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	60
5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	55





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Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

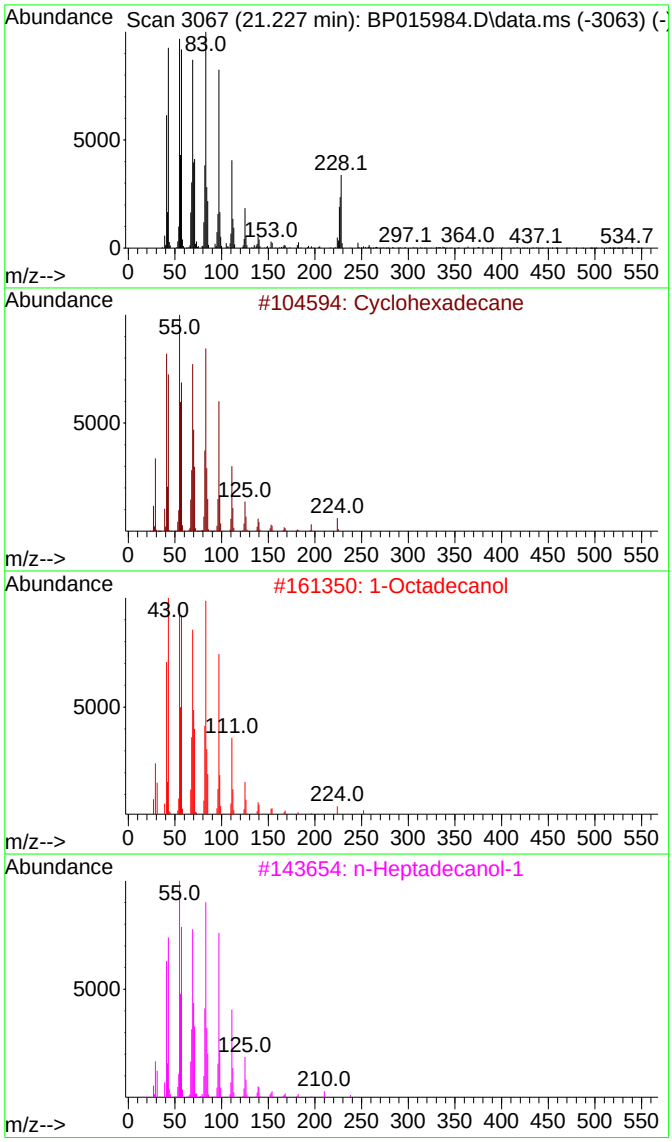
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

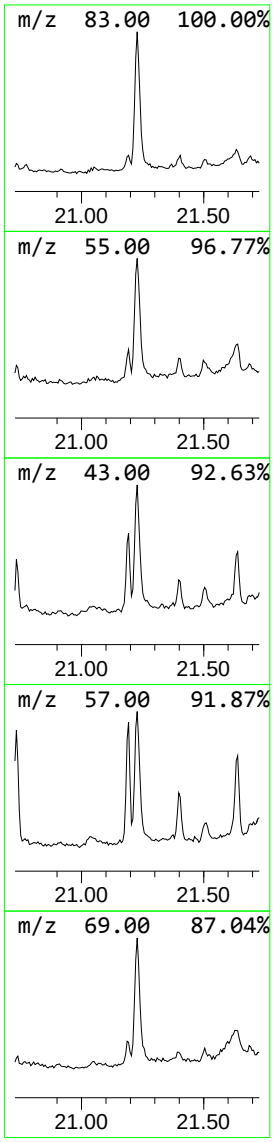
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic21.227 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.227	3.76 ng/ul	1271700	Chrysene-d12	21.545

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	94
2	1-Octadecanol	270	C18H38O	000112-92-5	91
3	n-Heptadecanol-1	256	C17H36O	001454-85-9	91
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	1-Octadecanol	270	C18H38O	000112-92-5	91





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Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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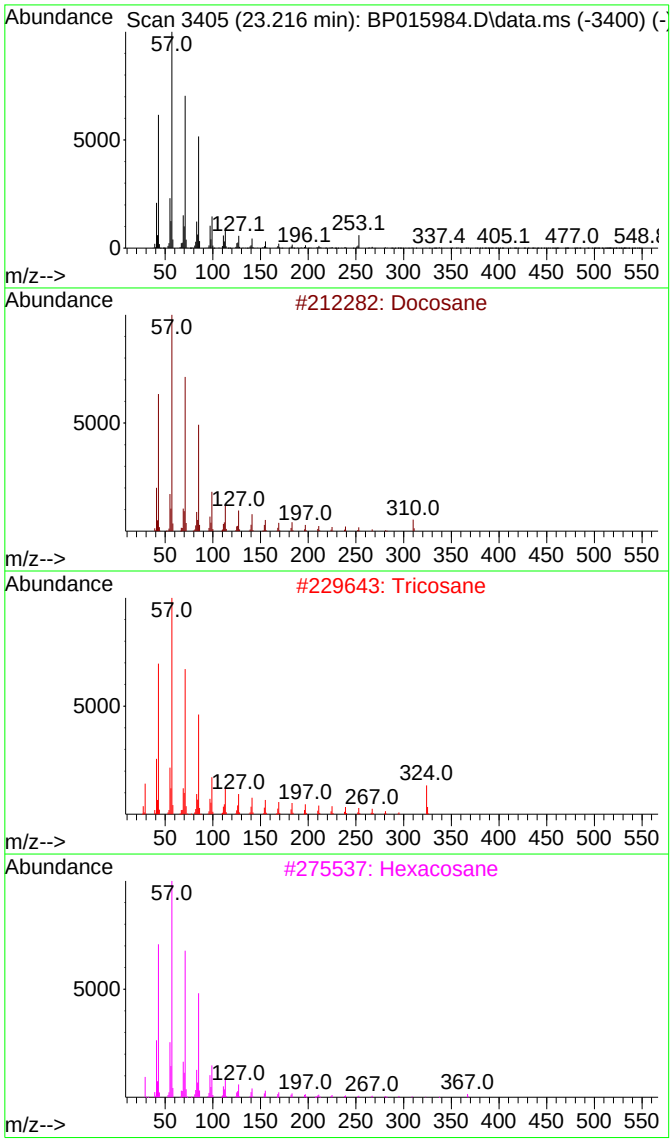
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

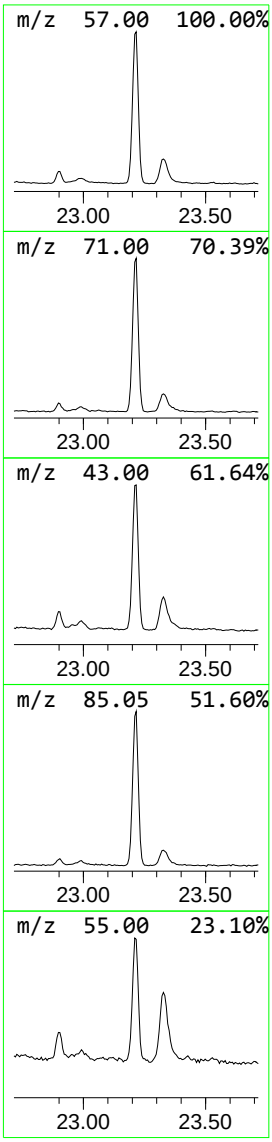
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.216	13.50 ng/ul	3359860	Perylene-d12	24.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane	310	C22H46	000629-97-0	95
2	Tricosane	324	C23H48	000638-67-5	93
3	Hexacosane	366	C26H54	000630-01-3	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

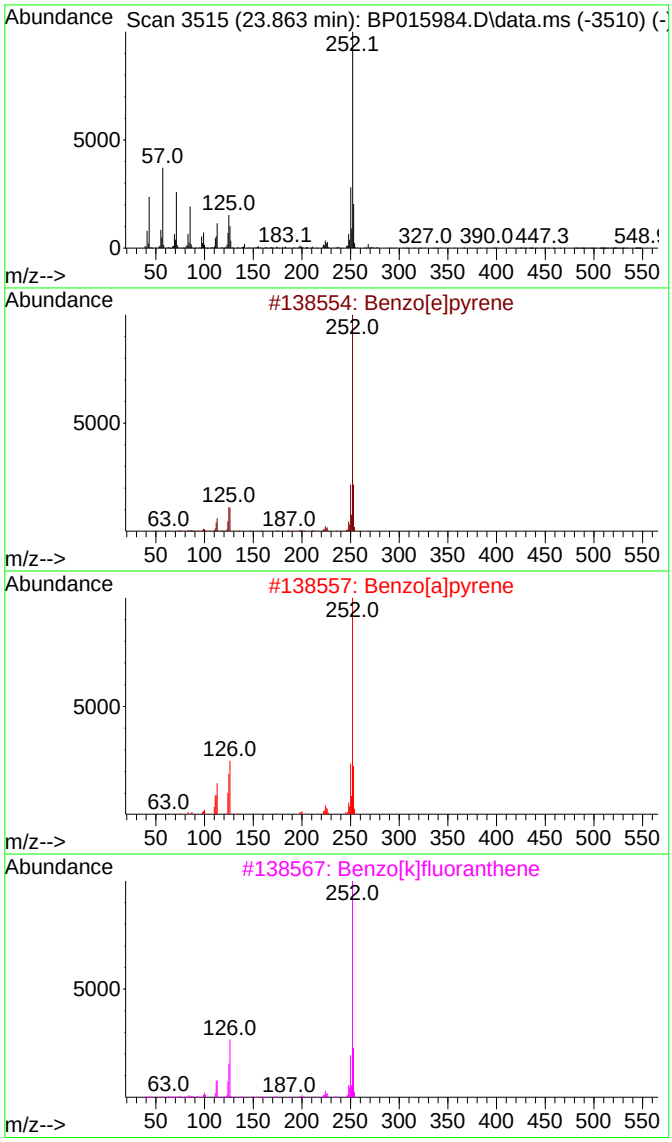
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

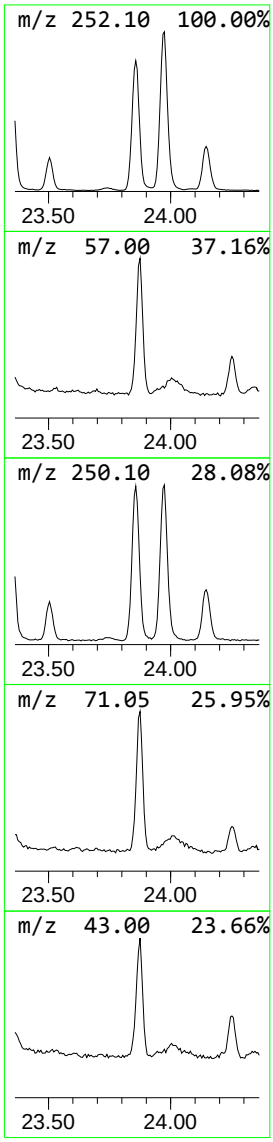
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.863	7.03 ng/ul	1749820	Perylene-d12	24.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[a]pyrene	252	C20H12	000050-32-8	95
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	92
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	92
5	Benzo[b]fluoranthene	252	C20H12	000205-99-2	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

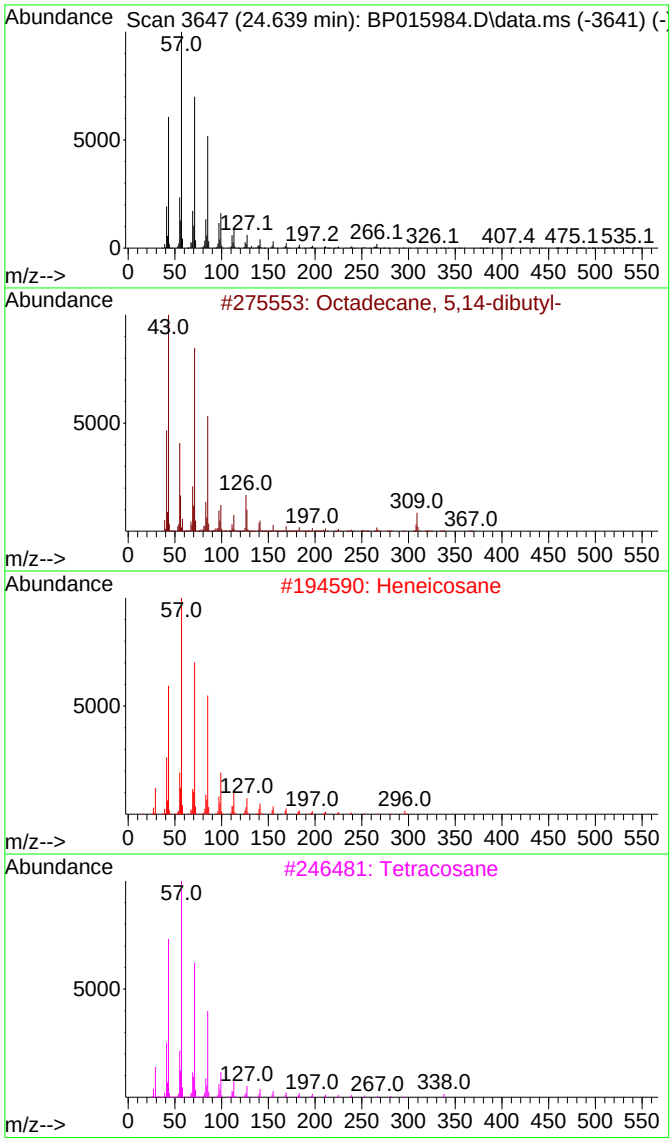
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

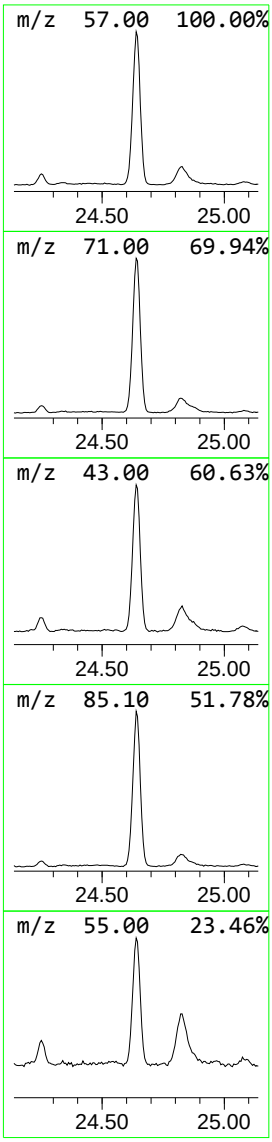
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.639	17.85 ng/ul	4442350	Perylene-d12	24.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Tetracosane	338	C24H50	000646-31-1	91
4	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5	Octacosane	394	C28H58	000630-02-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015984.D
Acq On : 04 Jul 2023 19:40
Operator : MA/JU
Sample : 03245-16
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGH4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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-Pentanone, 4-...	5.105	2.8	ng/ul	494340	1	8.040	3476180	20.0
Aceto phenone	8.899	2.6	ng/ul	457719	1	8.040	3476180	20.0
Benzophenone	16.057	4.4	ng/ul	1398830	3	14.693	6371390	20.0
n-Hexadecanoic ...	18.310	11.5	ng/ul	3815540	4	17.451	6624060	20.0
Octadecanoic acid	19.533	2.9	ng/ul	979776	5	21.545	6760770	20.0
Benzo[b]naphtho...	21.192	2.7	ng/ul	899425	5	21.545	6760770	20.0
(DEL) Alkane: C...	21.227	3.8	ng/ul	1271700	5	21.545	6760770	20.0
(DEL) Alkane: S...	23.216	13.5	ng/ul	3359860	6	24.086	4976460	20.0
Benzo[e]pyrene	23.863	7.0	ng/ul	1749820	6	24.086	4976460	20.0
(DEL) Alkane: S...	24.639	17.9	ng/ul	4442350	6	24.086	4976460	20.0