LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_M\Methods\8270-BM080122.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM036202.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.716	271	277	291	rBV2	45167	175229	1.64%	0.182%
2	4.940	310	315	319	rVB	78230	108626	1.02%	0.113%
3	5.416	389	396	411	rBV	3777807	5431750	50.79%	5.639%
4	7.034	663	671	684	rBV	3486525	5564957	52.03%	5.777%
5	7.828	797	806	815	rBV	908142	1430461	13.37%	1.485%
6	9.004	998	1006	1021	rBV	2126991	3568304	33.36%	3.704%
7	10.633	1273	1283	1289	rBV	1209068	1990959	18.61%	2.067%
8	13.098	1694	1702	1717	rBV	5359599	7758803	72.54%	8.054%
9	14.198	1884	1889	1898	rVB	98517	151066	1.41%	0.157%
10	14.363	1910	1917	1925	rBV2	52984	140925	1.32%	0.146%
11	14.474	1928	1936	1942	rVV	1720234	2532875	23.68%	2.629%
12	14.533	1942	1946	1954	rVB	166700	268428	2.51%	0.279%
13	14.874	1998	2004	2012	rBV	117575	196481	1.84%	0.204%
14	15.521	2108	2114	2125	rVB2	227880	389372	3.64%	0.404%
15	15.815	2159	2164	2171	rVB	68394	108838	1.02%	0.113%
16	15.962	2178	2189	2198	rBV	4320885	6036171	56.44%	6.266%
17	16.521	2273	2284	2289	rBV3	73287	180551	1.69%	0.187%
18	16.621	2296	2301	2307	rBV3	101317	178632	1.67%	0.185%
19	17.033	2367	2371	2379	rVB	128823	208285	1.95%	0.216%
20	17.115	2381	2385	2392	rBV2	80577	126689	1.18%	0.132%
21	17.215	2392	2402	2405	rBV	2103608	2941920	27.51%	3.054%
22	17.256	2405	2409	2415	rVB	2390518	3102104	29.00%	3.220%
23	17.345	2420	2424	2429	rVV	536289	766193	7.16%	0.795%
24	17.398	2429	2433	2445	rVB8	40991	132895	1.24%	0.138%
25	17.604	2459	2468	2470	rBV3	161069	294966	2.76%	0.306%
26	17.621	2470	2471	2483	rVB	178157	240432	2.25%	0.250%
27	17.886	2512	2516	2521	rVB	174335	212087	1.98%	0.220%
28	18.092	2542	2551	2552	rBV	259747	357855	3.35%	0.371%
29	18.115	2552	2555	2557	rBV	379603	404490	3.78%	0.420%
30	18.221	2569	2573	2578	rVB	133876	178377	1.67%	0.185%
31	18.286	2578	2584	2588	rBV3	407327	707755	6.62%	0.735%
32	18.321	2588	2590	2597	rVB	177404	209931	1.96%	0.218%
33	18.403	2599	2604	2608	rBV3	102547	172277	1.61%	0.179%
34	18.592	2631	2636	2640	rBV	191582	257461	2.41%	0.267%
35	18.639	2640	2644	2650	rVB	92057	129655	1.21%	0.135%
36	18.974	2696	2701	2705	rBV	529264	785622	7.35%	0.816%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_M\Methods\8270-BM080122.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	19.062	2712	2716	2720	rVB3	260317	309095	2.89%	0.321%
38	19.150	2726	2731	2735	rBV3	121453	233296	2.18%	0.242%
39	19.192	2735	2738	2741	rVV	207420	285473	2.67%	0.296%
40	19.227	2741	2744	2747	rVV	544574	635563	5.94%	0.660%
41	19.274	2747	2752	2759	rVB	3370362	4740988	44.33%	4.922%
42	19.603	2803	2808	2809	rBV	521119	643476	6.02%	0.668%
43	19.633	2809	2813	2822	rVV	2780078	3790317	35.44%	3.935%
44	19.703	2822	2825	2830	rVV3	124017	171959	1.61%	0.179%
45	19.839	2840	2848	2853	rBV	9011151	10695477	100.00%	11.103%
46	19.956	2863	2868	2875	rBV3	187908	457166	4.27%	0.475%
47	20.092	2886	2891	2893	rBV2	147206	261947	2.45%	0.272%
48	20.127	2893	2897	2904	rVB	586745	895087	8.37%	0.929%
49	20.227	2910	2914	2918	rBV	405539	530712	4.96%	0.551%
50	20.286	2919	2924	2928	rVV2	269532	453649	4.24%	0.471%
51	20.321	2928	2930	2934	rVB	88130	112025	1.05%	0.116%
52	20.427	2944	2948	2951	rBV2	156528	219322	2.05%	0.228%
53	20.480	2953	2957	2965	rVB2	276648	468211	4.38%	0.486%
54	20.603	2975	2978	2981	rVB	209175	201301	1.88%	0.209%
55	20.874	3021	3024	3030	rVB	165377	244075	2.28%	0.253%
56	21.039	3049	3052	3056	rBV	215855	256153	2.39%	0.266%
57	21.080	3056	3059	3060	rBV	244657	234497	2.19%	0.243%
58	21.109	3060	3064	3071	rVB4	1196391	1713713	16.02%	1.779%
59	21.391	3105	3112	3116	rBV2	3444639	6487098	60.65%	6.734%
60	21.427	3116	3118	3125	rVB	1573684	1914591	17.90%	1.988%
61	21.544	3135	3138	3144	rBV	335427	495290	4.63%	0.514%
62	22.015	3216	3218	3224	rVB3	422819	549250	5.14%	0.570%
63	23.021	3384	3389	3395	rBV2	1699740	3800776	35.54%	3.946%
64	23.533	3471	3476	3485	rVB	748065	1459871	13.65%	1.515%
65	23.633	3488	3493	3502	rVB	1243841	2274587	21.27%	2.361%
66	23.738	3505	3511	3517	rBV2	1814371	3435419	32.12%	3.566%
67	24.327	3608	3611	3622	rVB2	467734	919612	8.60%	0.955%

Sum of corrected areas: 96331418

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
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Acq On : 10 Aug 2022 21:13
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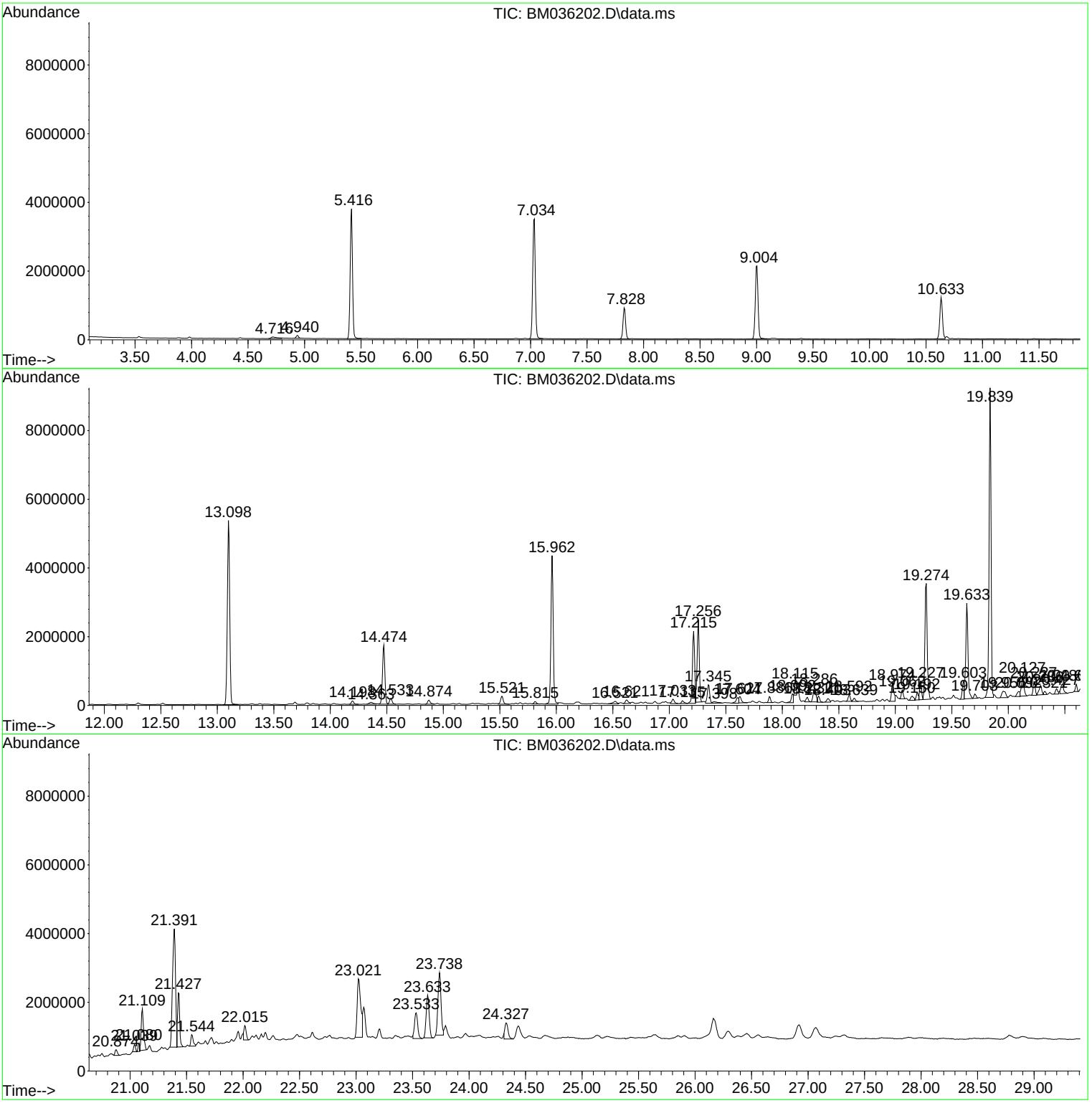
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

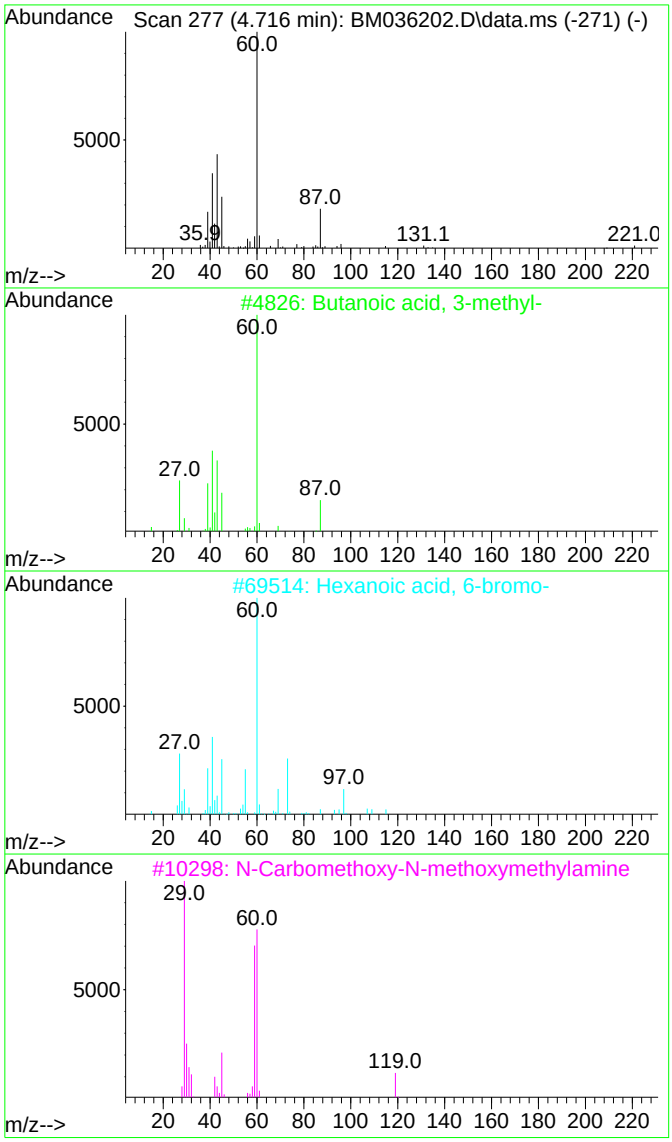
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

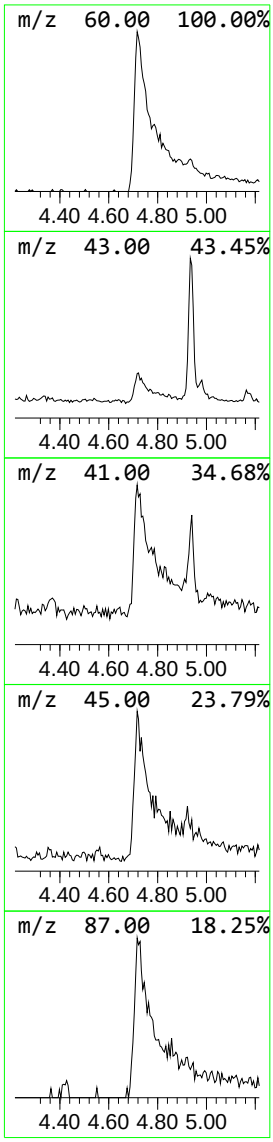
****

Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.716	2.45 ng	175229	1,4-Dichlorobenzene-d4	7.828

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
2	Hexanoic acid, 6-bromo-	194	C6H11BrO2	004224-70-8	40
3	N-Carbomethoxy-N-methoxymethylamine	119	C4H9NO3	153654-07-0	39
4	Hexanoic acid	116	C6H12O2	000142-62-1	39
5	9-Fluorenone, thiosemicarbazone	253	C14H11N3S	1010293-93-5	36





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

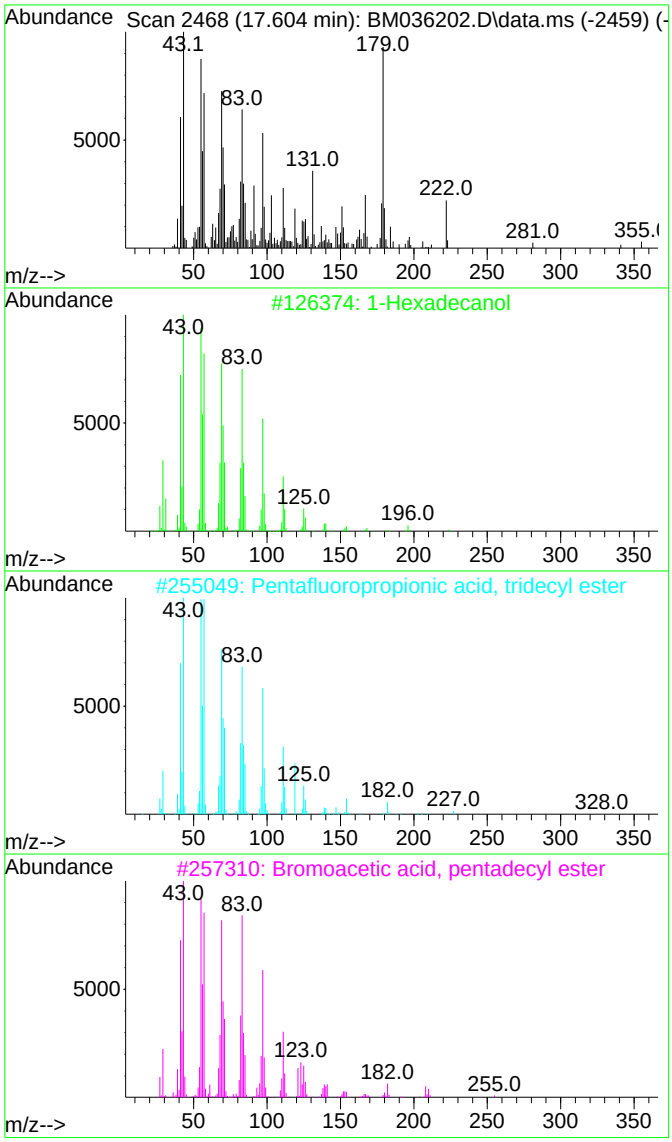
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

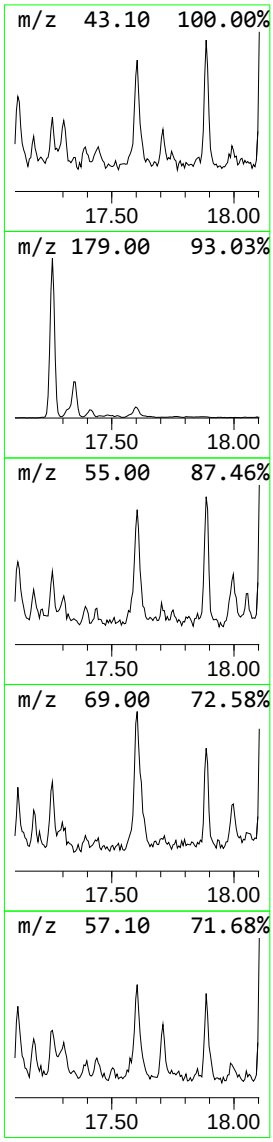
****

Peak Number 3 1-Hexadecanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.604	2.01 ng	294966	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	55
2	Pentafluoropropionic acid, tride...	346	C16H27F5O2	959261-22-4	45
3	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	43
4	Phenanthridine	179	C13H9N	000229-87-8	42
5	Pentafluoropropionic acid, undec...	318	C14H23F5O2	959014-11-0	41





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

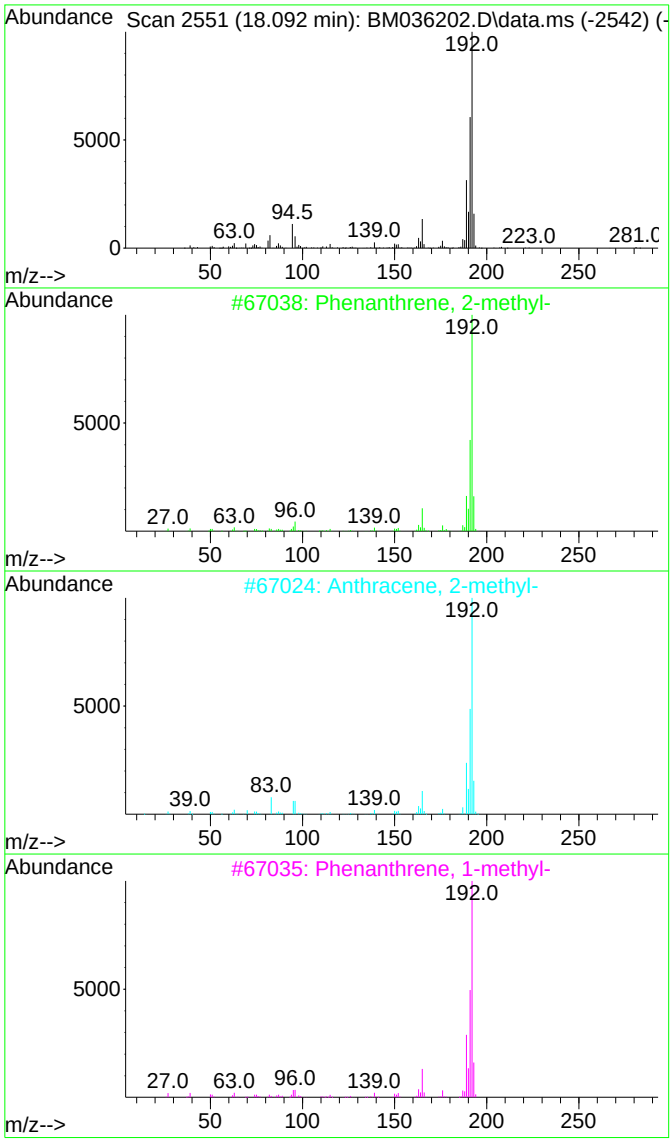
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

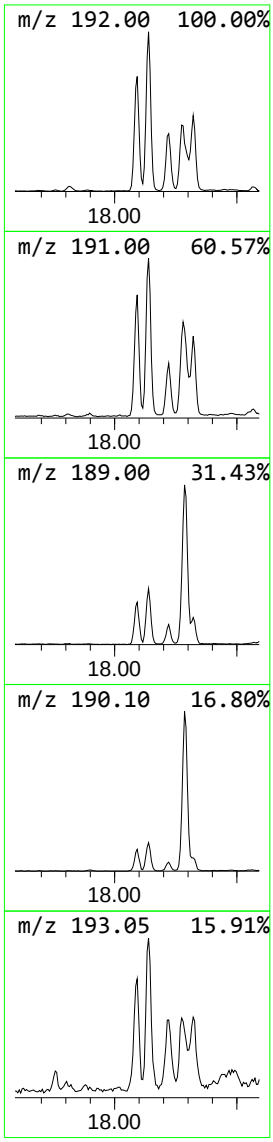
****

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	2.43 ng	357855	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

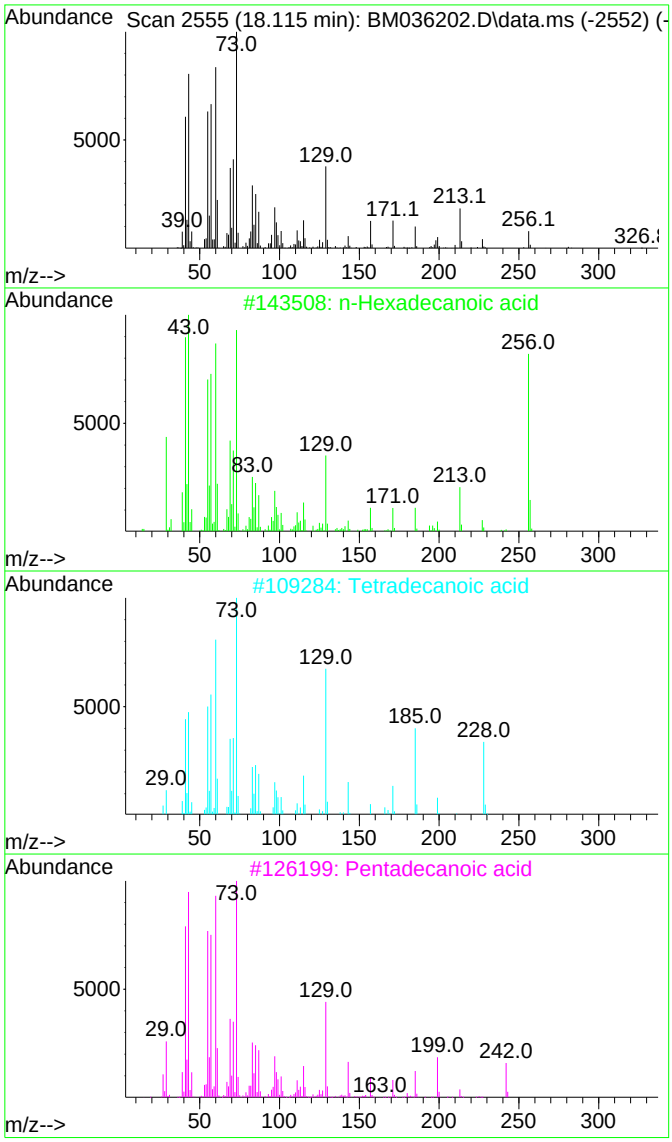
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

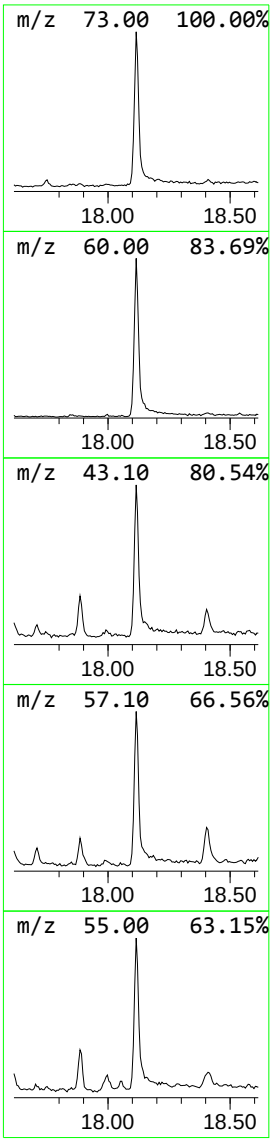
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.115	2.75 ng	404490	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tridecanoic acid	214	C13H26O2	000638-53-9	83
5	8-Methylnonanoic acid	172	C10H20O2	005963-14-4	49





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

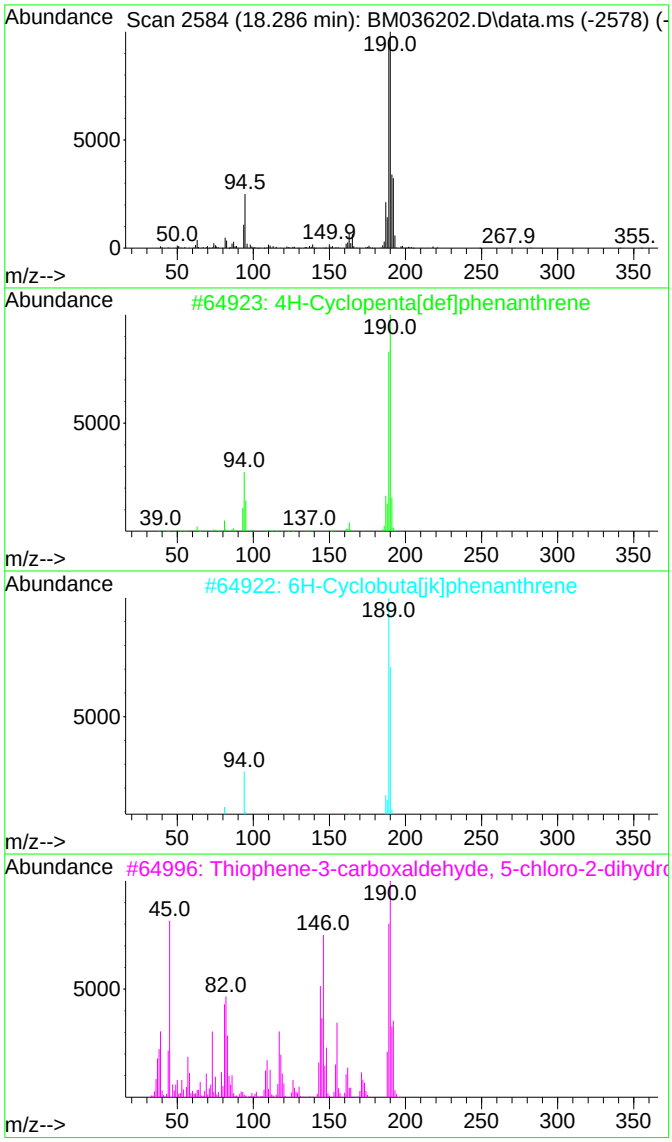
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

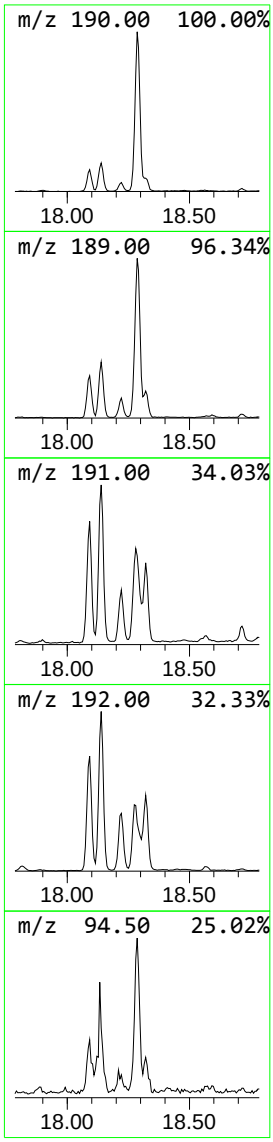
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	4.81 ng	707755	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	33
5	1-Aminocyclopentanecarboxylic ac...	361	C18H32ClNO4	1000329-17-0	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
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Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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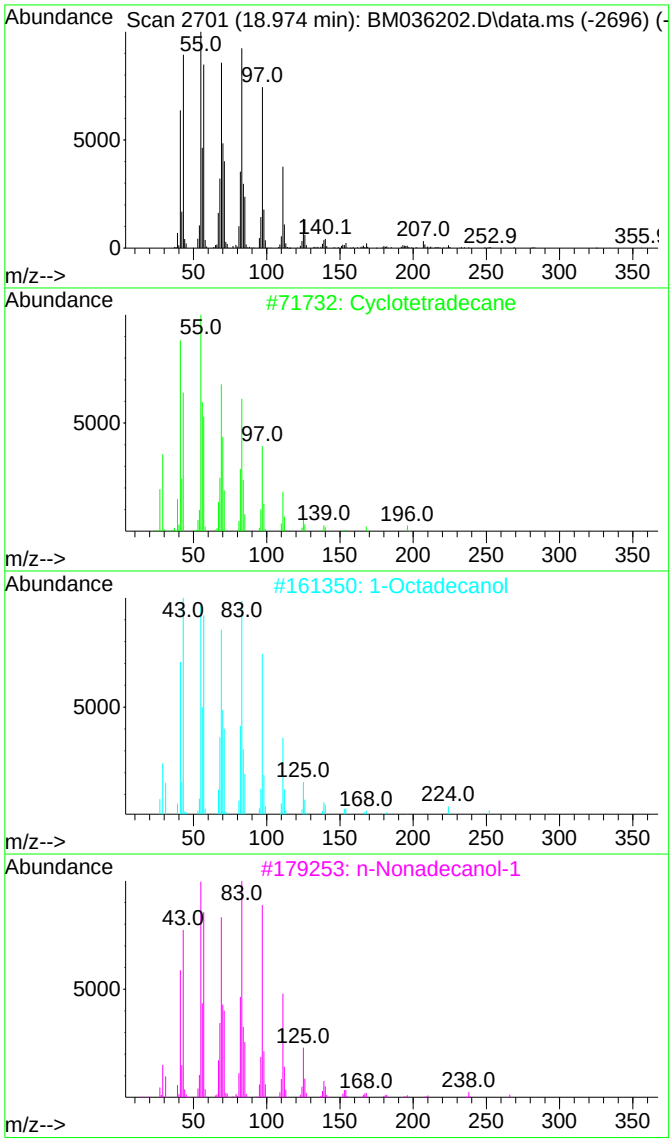
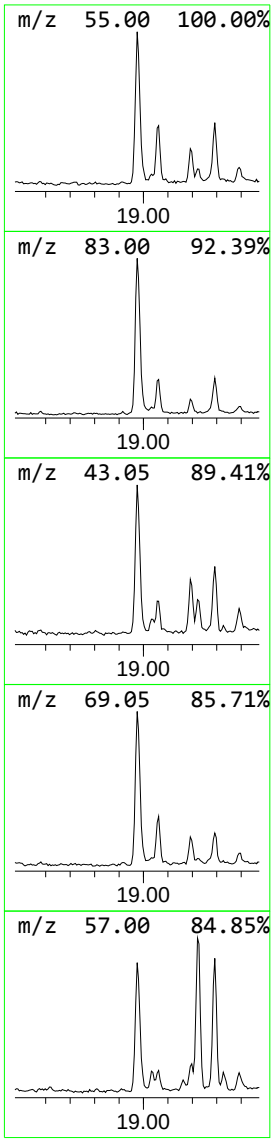
Peak Number 7 Cyclotetradecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.974	5.34 ng	785622	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	97
2	1-Octadecanol	270	C18H38O	000112-92-5	94
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4	1-Hexadecanol	242	C16H34O	036653-82-4	93
5	n-Heptadecanol-1	256	C17H36O	001454-85-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

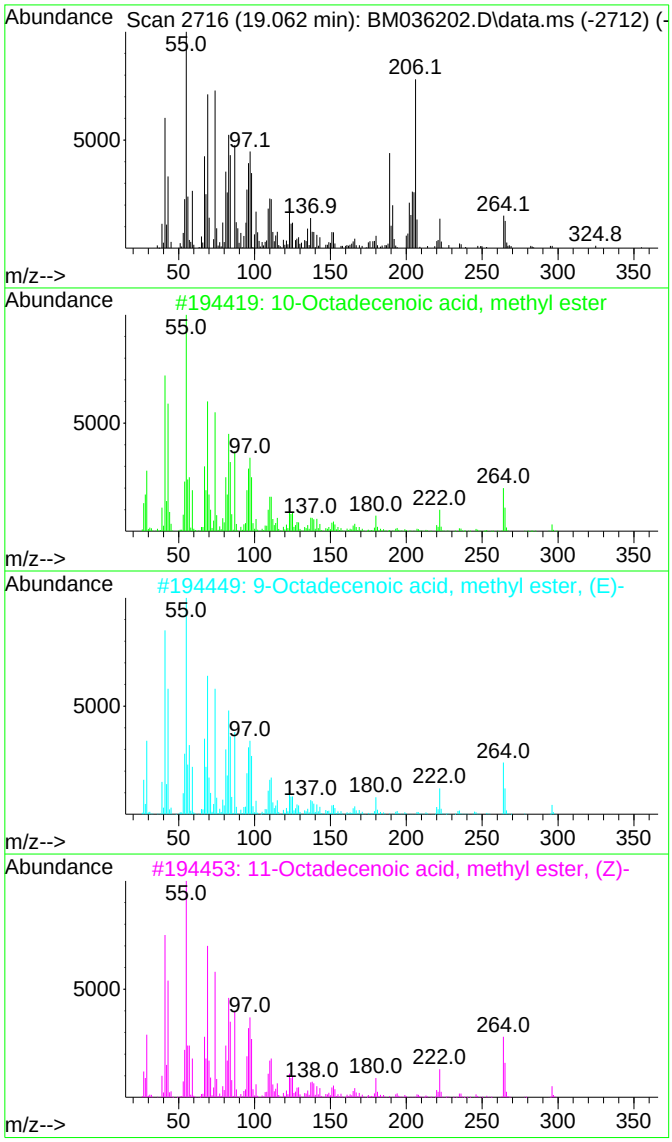
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

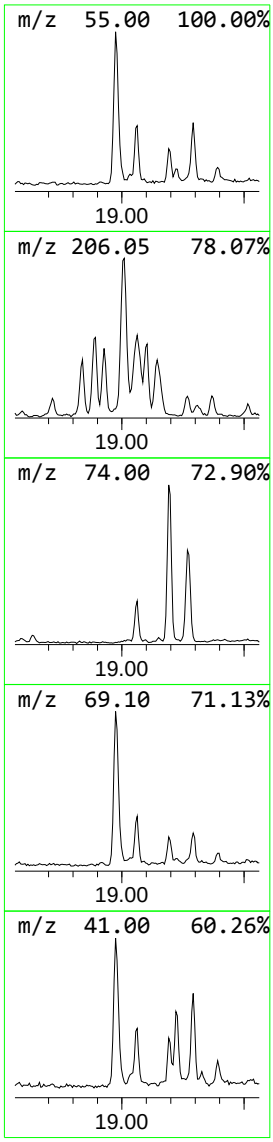
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 10-Octadecenoic acid, methy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.062	2.10 ng	309095	Phenanthrene-d10	17.215

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	97
2	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	97
3	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	97
4	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	97
5	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	96





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Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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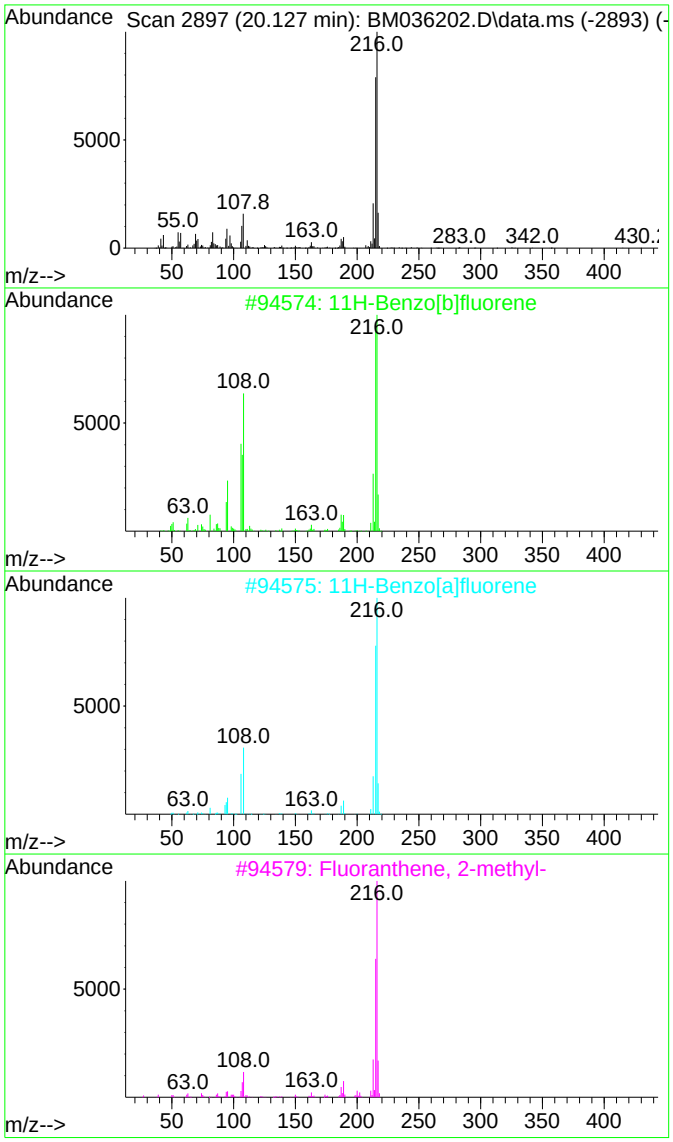
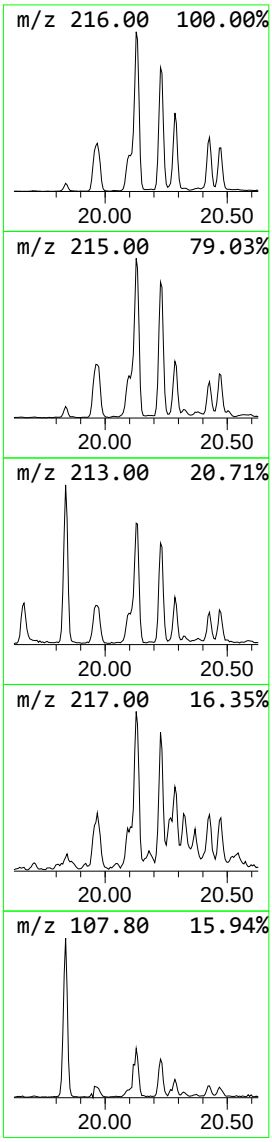
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.127	2.76 ng	895087	Chrysene-d12	21.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81





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Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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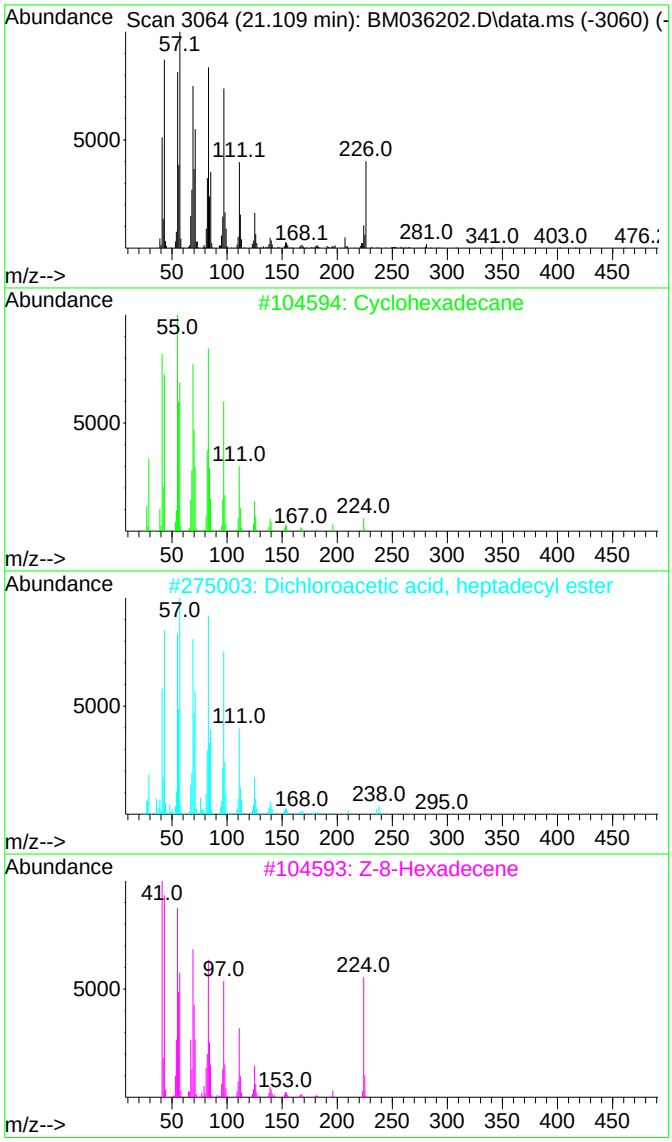
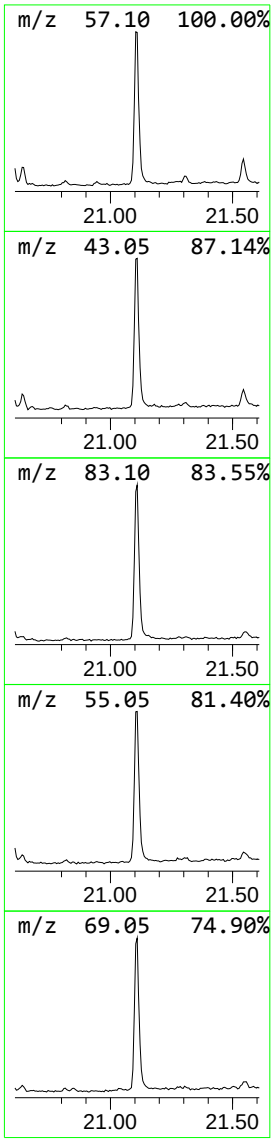
Peak Number 10 Cyclohexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.109	5.28 ng	1713710	Chrysene-d12	21.392

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	94
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
3	Z-8-Hexadecene	224	C16H32	1000130-87-5	91
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	90





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Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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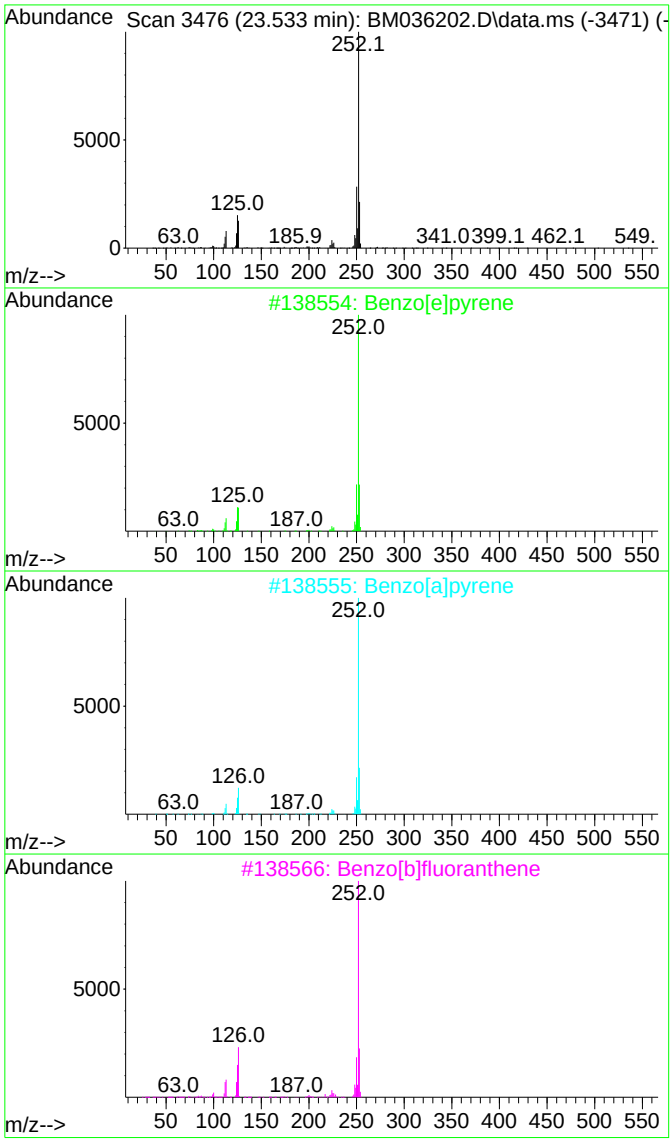
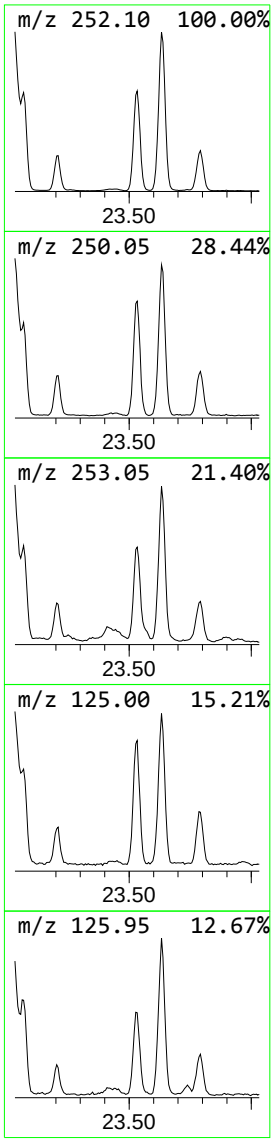
Peak Number 11 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.532	8.50 ng	1459870	Perylene-d12	23.738

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	96
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4	Perylene	252	C20H12	000198-55-0	94
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	93





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Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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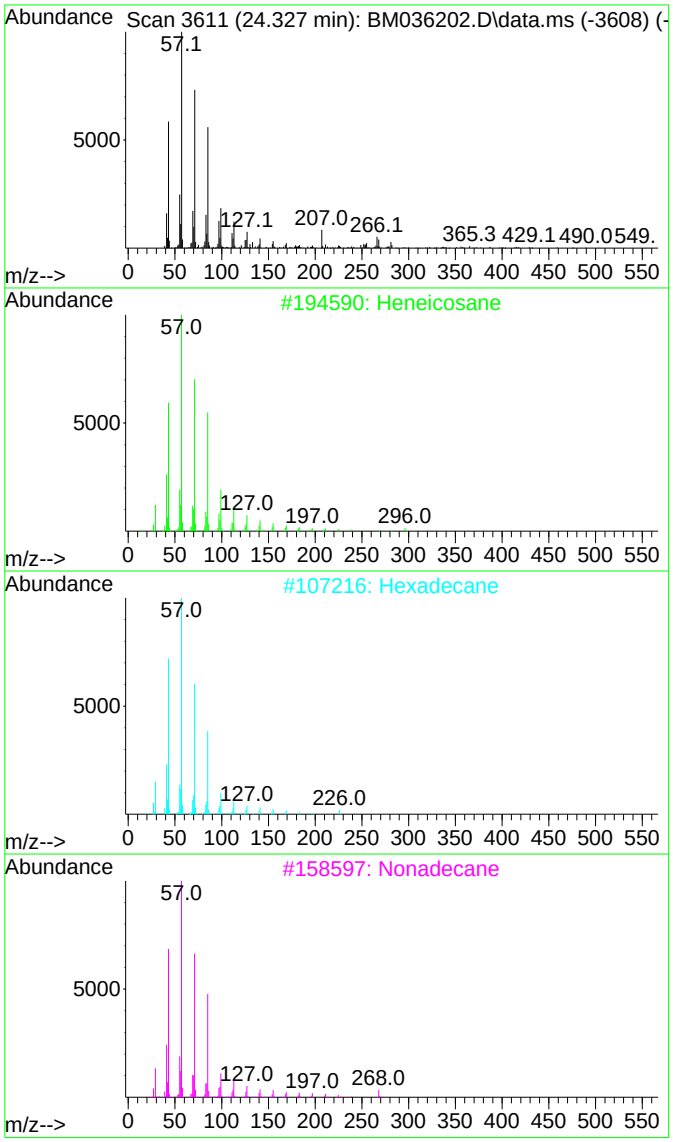
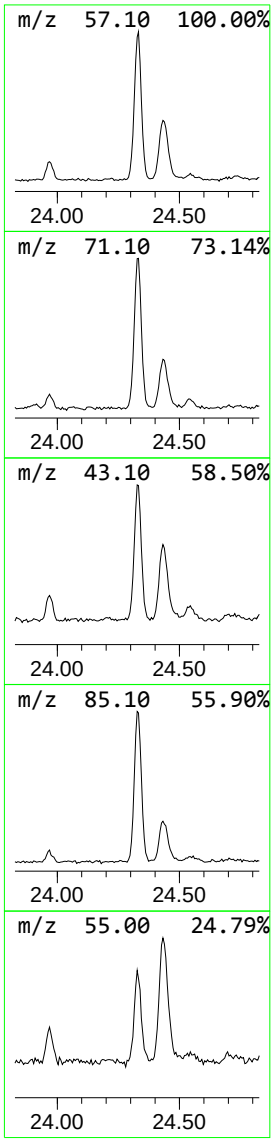
Peak Number 12 Heneicosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.327	5.35 ng	919612	Perylene-d12	23.738

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	96
2	Hexadecane	226	C16H34	000544-76-3	94
3	Nonadecane	268	C19H40	000629-92-5	94
4	Docosane	310	C22H46	000629-97-0	87
5	Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	86





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
Data File : BM036202.D
Acq On : 10 Aug 2022 21:13
Operator : CG/JU
Sample : N4086-01
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
RLA-4711-WC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butanoic acid, ...	4.716	2.5	ng	175229	1	7.828	1430460	20.0
1-Hexadecanol	17.604	2.0	ng	294966	4	17.215	2941920	20.0
Phenanthrene, 2...	18.092	2.4	ng	357855	4	17.215	2941920	20.0
n-Hexadecanoic ...	18.115	2.8	ng	404490	4	17.215	2941920	20.0
4H-Cyclopenta[d...	18.286	4.8	ng	707755	4	17.215	2941920	20.0
Cyclotetradecane	18.974	5.3	ng	785622	4	17.215	2941920	20.0
10-Octadecenoic...	19.062	2.1	ng	309095	4	17.215	2941920	20.0
11H-Benzo[b]flu...	20.127	2.8	ng	895087	5	21.392	6487100	20.0
Cyclohexadecane	21.109	5.3	ng	1713710	5	21.392	6487100	20.0
Benzo[e]pyrene	23.532	8.5	ng	1459870	6	23.738	3435420	20.0
Heneicosane	24.327	5.3	ng	919612	6	23.738	3435420	20.0