LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

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_P\METHODS\8270-BP112719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.617	595	600	613	rVB	245745	389549	15.79%	1.425%
2	6.211	695	701	710	rBV	1517149	1874454	76.00%	6.857%
3	7.752	948	963	974	rBV	1693479	2223620	90.16%	8.135%
4	7.940	989	995	1004	rBV	1651767	1988848	80.64%	7.276%
5	8.305	1051	1057	1063	rBV	443758	497338	20.17%	1.819%
6	8.546	1092	1098	1103	rBV	1268728	1480395	60.02%	5.416%
7	9.181	1200	1206	1211	rBV	1097617	1308412	53.05%	4.787%
8	10.334	1397	1402	1407	rBV	519459	614251	24.91%	2.247%
9	10.969	1504	1510	1519	rBV2	22781	38386	1.56%	0.140%
10	12.116	1696	1705	1709	rBV	1887903	2170536	88.01%	7.941%
11	12.781	1813	1818	1823	rBV	192098	216747	8.79%	0.793%
12	13.198	1883	1889	1894	rBV	627458	759155	30.78%	2.777%
13	14.093	2037	2041	2046	rVB	23902	26633	1.08%	0.097%
14	14.487	2102	2108	2121	rBV	1423859	1757614	71.26%	6.430%
15	15.592	2291	2296	2299	rBV	724795	839956	34.06%	3.073%
16	15.628	2299	2302	2307	rVB	334540	354821	14.39%	1.298%
17	15.704	2312	2315	2318	rBV	59121	58527	2.37%	0.214%
18	15.951	2353	2357	2361	rVB	30499	32672	1.32%	0.120%
19	16.345	2420	2424	2428	rBV	34378	40671	1.65%	0.149%
20	16.387	2428	2431	2437	rVB	49420	58654	2.38%	0.215%
21	16.481	2444	2447	2454	rBV4	152920	250058	10.14%	0.915%
22	16.534	2454	2456	2462	rVB	28770	35894	1.46%	0.131%
23	16.792	2495	2500	2507	rVB2	40372	68836	2.79%	0.252%
24	17.134	2552	2558	2561	rBV2	17962	29459	1.19%	0.108%
25	17.228	2568	2574	2579	rBV3	29564	55444	2.25%	0.203%
26	17.339	2587	2593	2597	rBV	564370	621264	25.19%	2.273%
27	17.569	2629	2632	2637	rVV2	38163	42134	1.71%	0.154%
28	17.639	2639	2644	2649	rVB	482543	605244	24.54%	2.214%
29	17.816	2670	2674	2677	rBV	29091	32252	1.31%	0.118%
30	17.875	2677	2684	2688	rVV	2264288	2466308	100.00%	9.023%
31	18.069	2711	2717	2719	rBV5	22979	42160	1.71%	0.154%
32	18.092	2719	2721	2725	rVB2	51469	55648	2.26%	0.204%
33	18.186	2734	2737	2740	rBV	26774	28451	1.15%	0.104%
34	18.228	2740	2744	2755	rVV4	57841	138651	5.62%	0.507%

LSC Area Percent Report

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Misc :
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Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

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_P\METHODS\8270-BP112719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	18.528	2792	2795	2804	rVB7	16359	29663	1.20%	0.109%
36	18.692	2820	2823	2828	rVV	58571	62009	2.51%	0.227%
37	18.757	2830	2834	2842	rVB2	51247	67437	2.73%	0.247%
38	18.904	2854	2859	2863	rBV2	48296	67399	2.73%	0.247%
39	18.945	2863	2866	2868	rBV	32002	34925	1.42%	0.128%
40	19.028	2877	2880	2888	rVB	57641	68127	2.76%	0.249%
41	19.110	2889	2894	2905	rBV3	305207	616561	25.00%	2.256%
42	19.233	2909	2915	2918	rVV2	834795	1189991	48.25%	4.353%
43	19.263	2918	2920	2930	rVB2	276105	355566	14.42%	1.301%
44	19.363	2934	2937	2942	rVB	31268	37960	1.54%	0.139%
45	19.528	2961	2965	2969	rVB	108031	118164	4.79%	0.432%
46	19.692	2989	2993	2996	rBV	17545	25759	1.04%	0.094%
47	19.733	2996	3000	3004	rVV	48389	64770	2.63%	0.237%
48	19.775	3004	3007	3012	rVB	30179	37139	1.51%	0.136%
49	19.916	3028	3031	3037	rVB	104725	123608	5.01%	0.452%
50	20.110	3060	3064	3069	rBV4	26454	44693	1.81%	0.164%
51	20.292	3091	3095	3100	rBV2	94654	119009	4.83%	0.435%
52	20.516	3128	3133	3136	rBV	241602	401799	16.29%	1.470%
53	20.639	3150	3154	3157	rVB	41401	45655	1.85%	0.167%
54	20.686	3157	3162	3165	rBV	111410	154485	6.26%	0.565%
55	20.733	3166	3170	3173	rBV6	31378	54990	2.23%	0.201%
56	20.857	3186	3191	3198	rVB	146942	207499	8.41%	0.759%
57	20.927	3198	3203	3210	rVB	195168	266045	10.79%	0.973%
58	21.004	3210	3216	3219	rBV	521204	724922	29.39%	2.652%
59	21.033	3219	3221	3225	rVB	53315	53808	2.18%	0.197%
60	21.122	3231	3236	3242	rVB	121238	188498	7.64%	0.690%
61	21.616	3314	3320	3326	rBV	120388	208176	8.44%	0.762%
62	21.680	3327	3331	3339	rVV7	22399	57304	2.32%	0.210%
63	22.186	3412	3417	3426	rVB2	58380	112101	4.55%	0.410%
64	22.445	3456	3461	3468	rVB3	23019	47798	1.94%	0.175%
65	22.622	3485	3491	3502	rVB3	94154	229431	9.30%	0.839%
66	22.839	3521	3528	3534	rBV5	28899	99500	4.03%	0.364%
67	23.104	3567	3573	3588	rVB2	79226	178455	7.24%	0.653%
68	23.357	3613	3616	3623	rVB2	20338	38392	1.56%	0.140%

Sum of corrected areas: 27334680

LSC Report - Integrated Chromatogram

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Operator : JU
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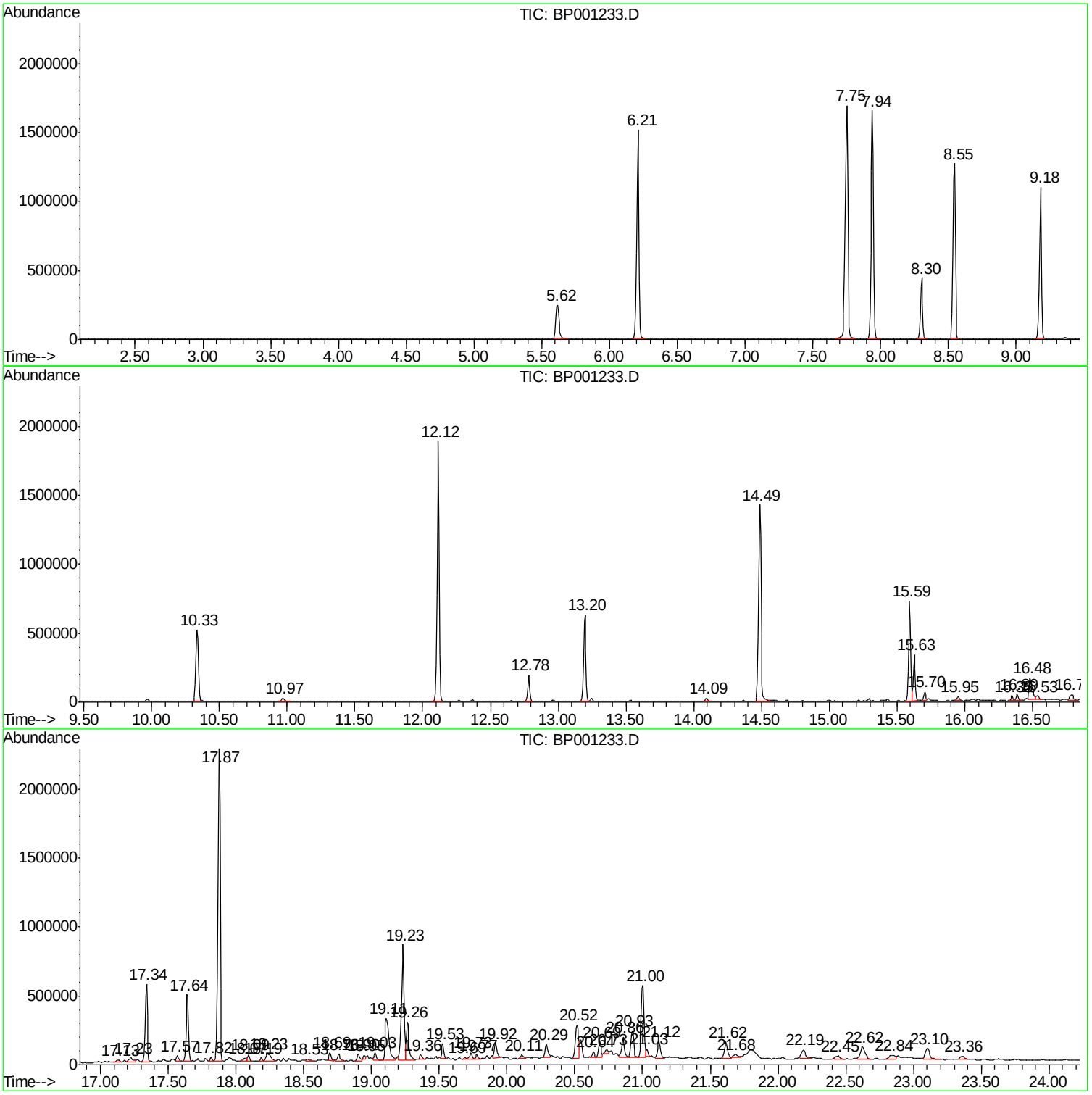
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001233.D
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Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

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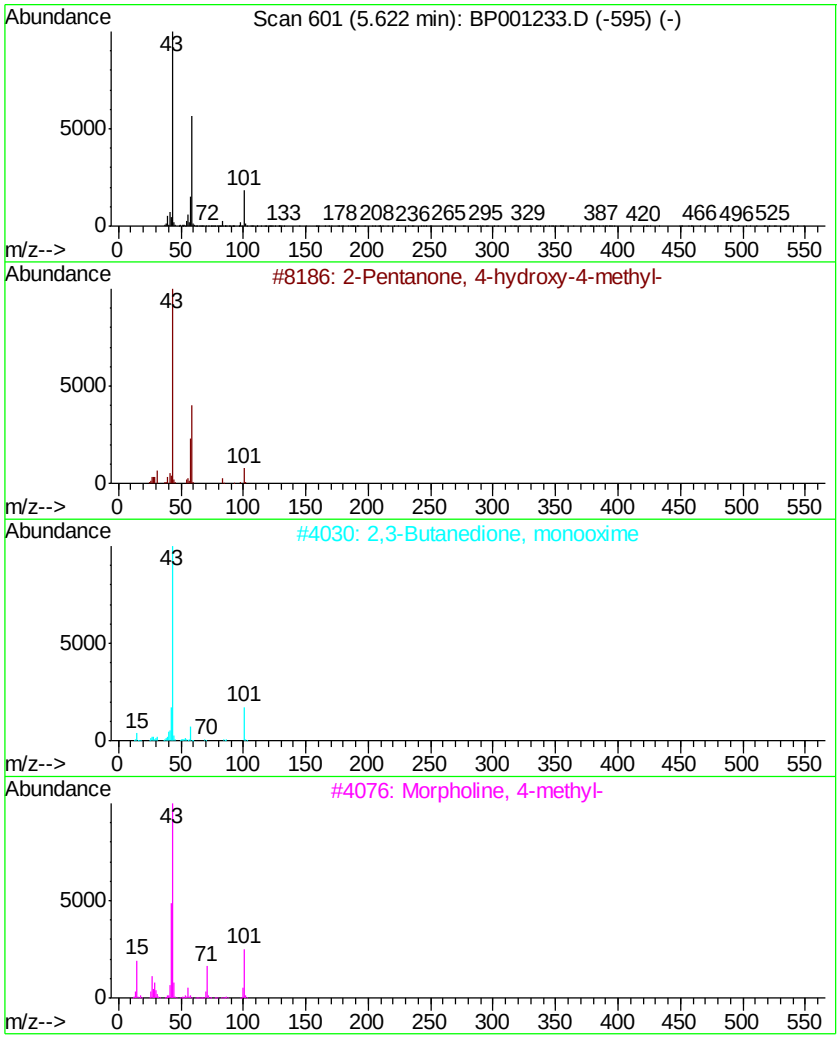
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

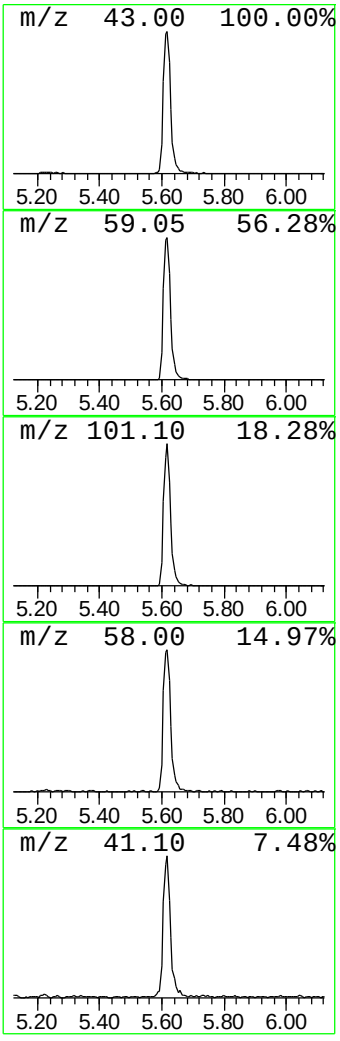
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.62	15.67 ng	389549	1,4-Dichlorobenzene-d4	8.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	Butane	58	C4H10	000106-97-8	9
5	Silane, trimethyl-	74	C3H10Si	000993-07-7	9





Library Search Compound Report

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Data File : BP001233.D
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Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

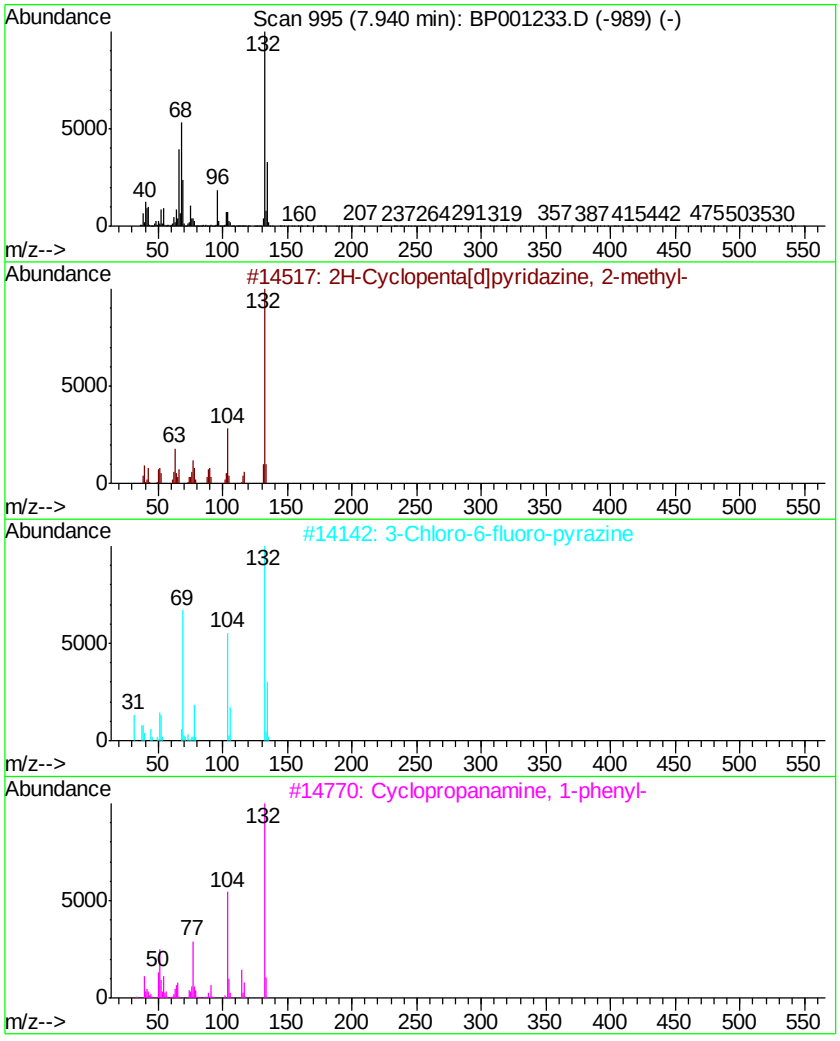
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

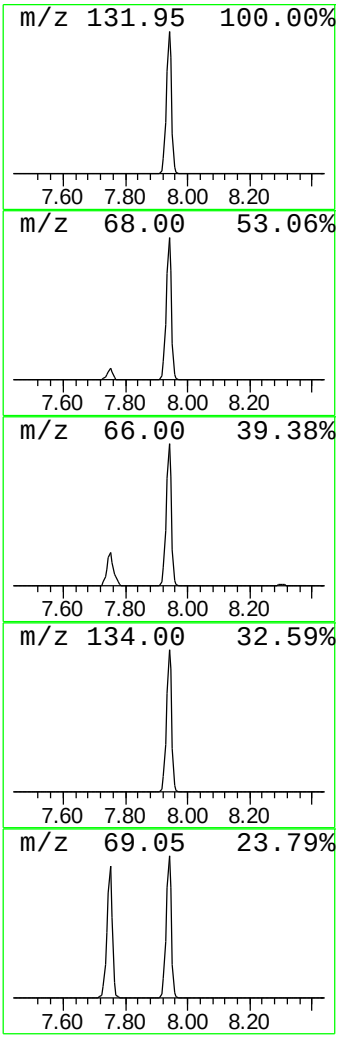
****

Peak Number 2 unknown7.94 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	79.98 ng	1988850	1,4-Dichlorobenzene-d4	8.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	14
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10
3	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10





Library Search Compound Report

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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
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TIC Library : C:\DATABASE\NIST11.L
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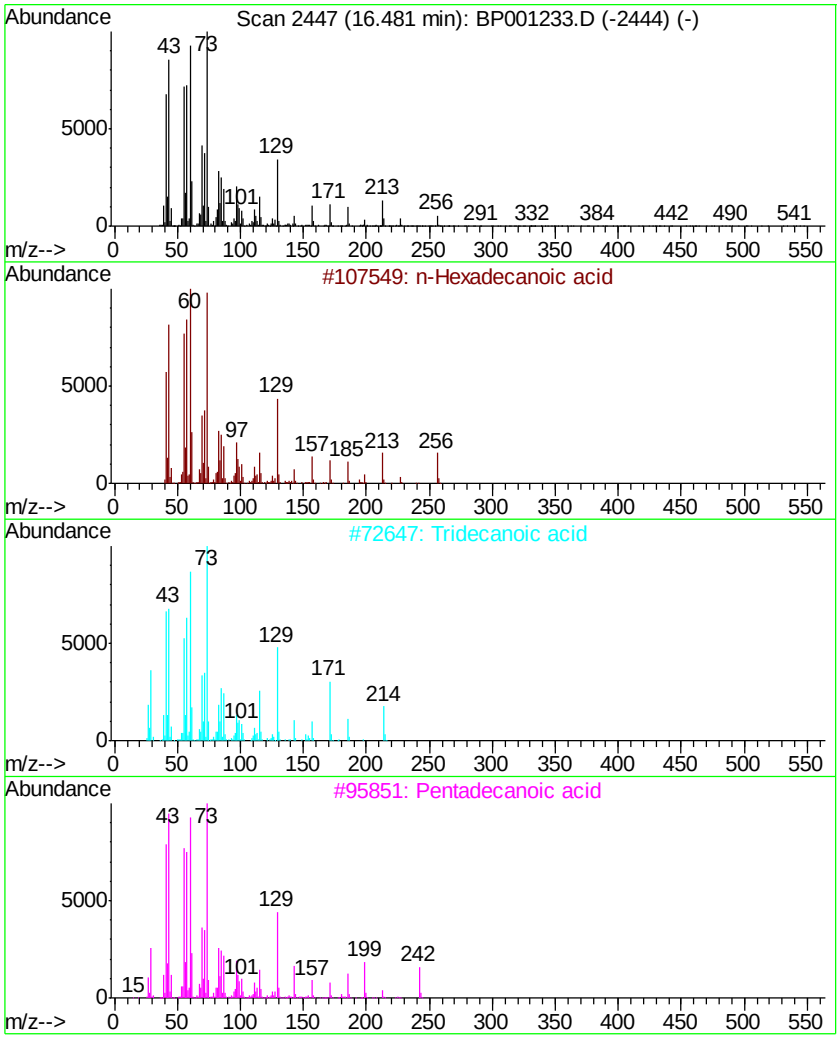
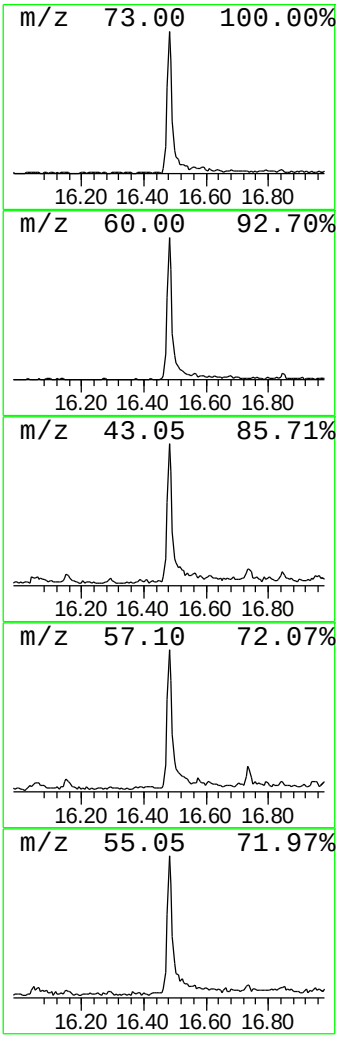
****

Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.48	5.95 ng	250058	Phenanthrene-d10	15.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5	Undecanoic acid	186	C11H22O2	000112-37-8	59





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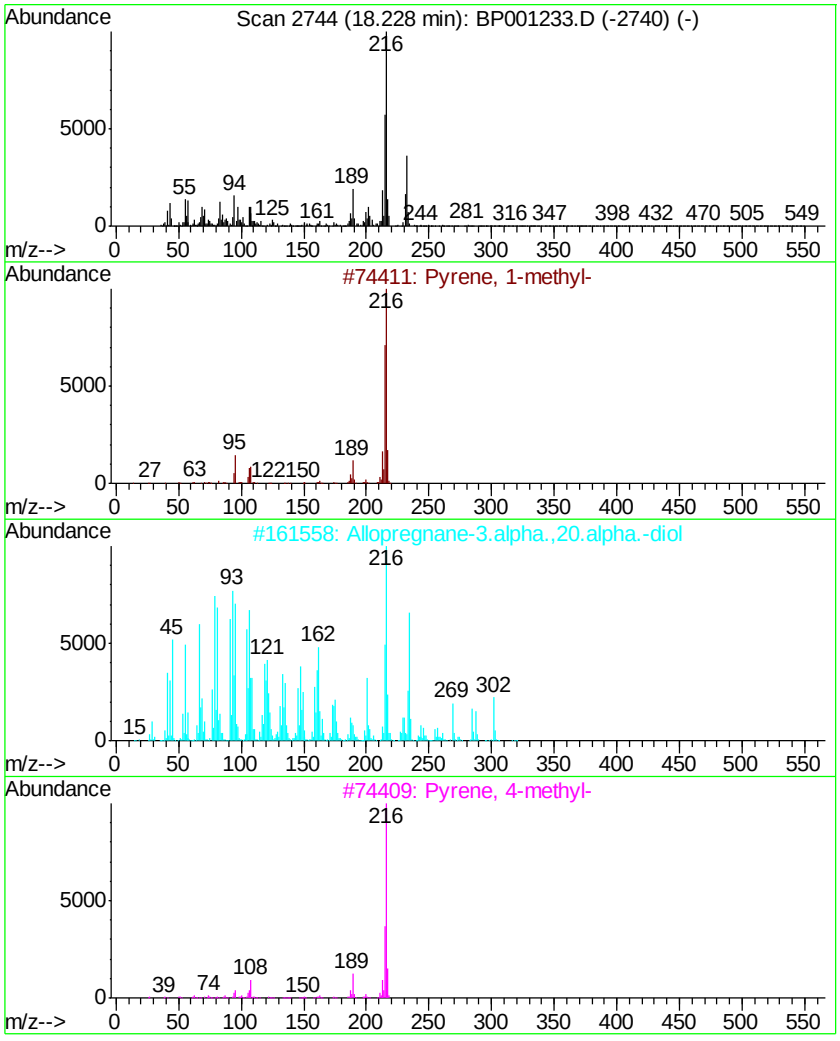
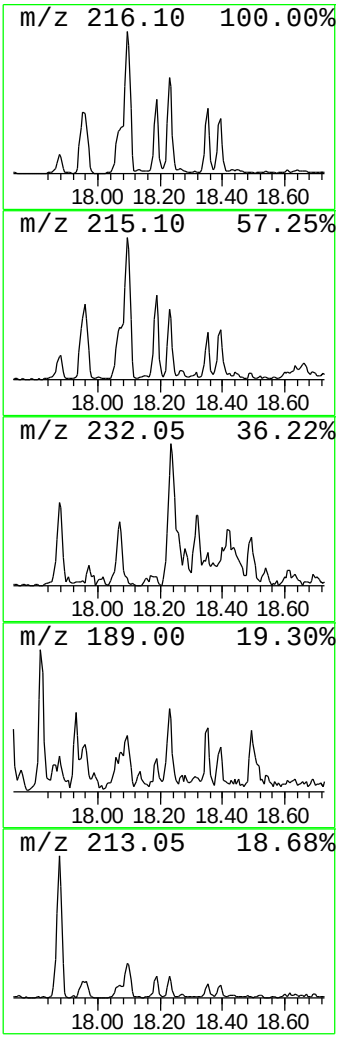
****

Peak Number 5 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.23	2.33 ng	138651	Chrysene-d12	19.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	Allopregnane-3.alpha.,20.alpha.-...	320	C21H36O2	000566-58-5	92
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	64
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	64





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Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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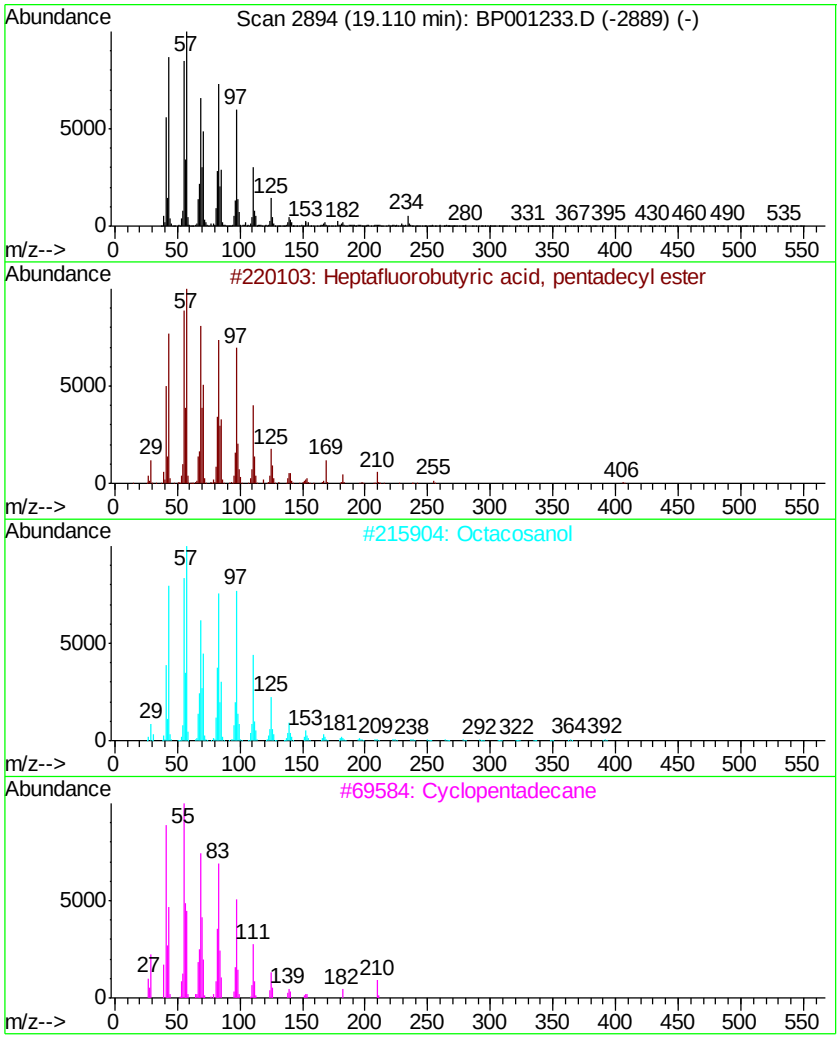
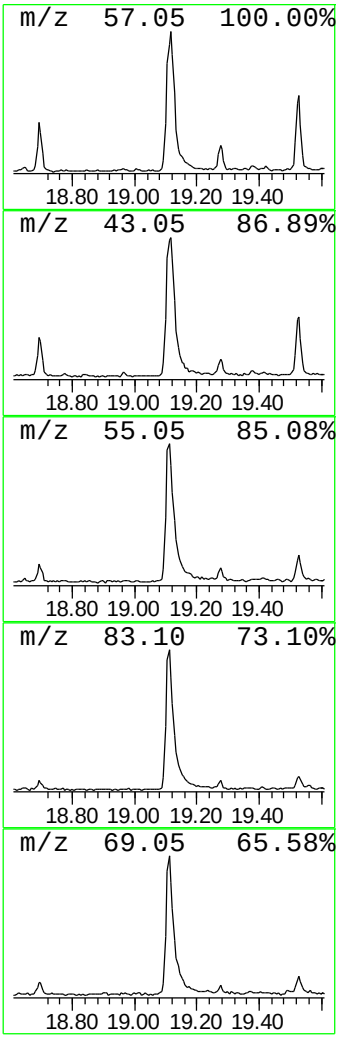
Peak Number 6 Heptafluorobutyric acid, pe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.11	10.36 ng	616561	Chrysene-d12	19.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2	Octacosanol	410	C28H58O	000557-61-9	94
3	Cyclopentadecane	210	C15H30	000295-48-7	92
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91





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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

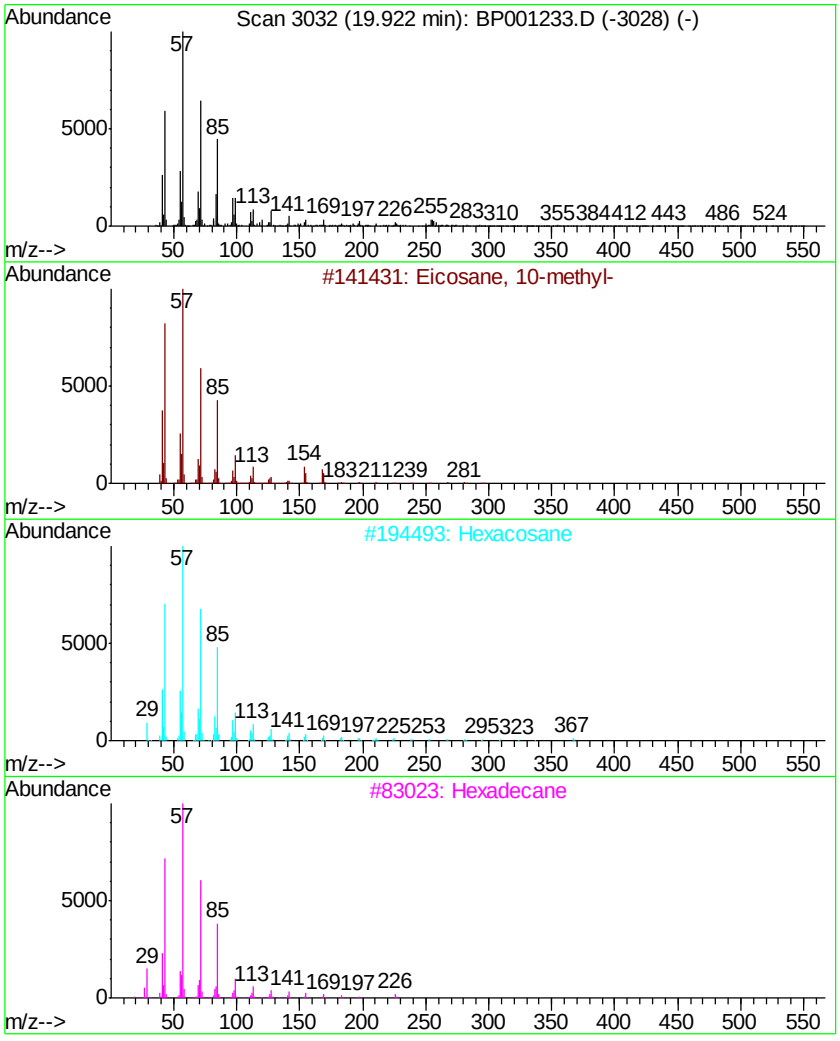
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

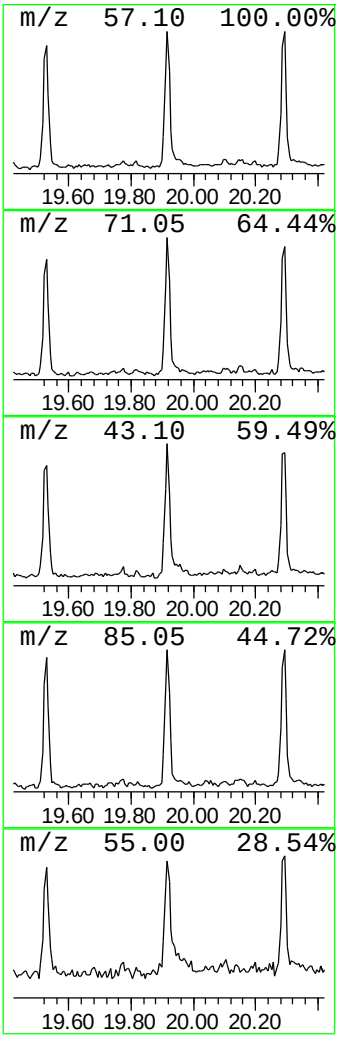
****

Peak Number 7 Eicosane, 10-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.92	2.08 ng	123608	Chrysene-d12	19.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2	Hexacosane	366	C26H54	000630-01-3	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Heptacosane	380	C27H56	000593-49-7	90
5	Octacosane	394	C28H58	000630-02-4	90





Library Search Compound Report

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Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

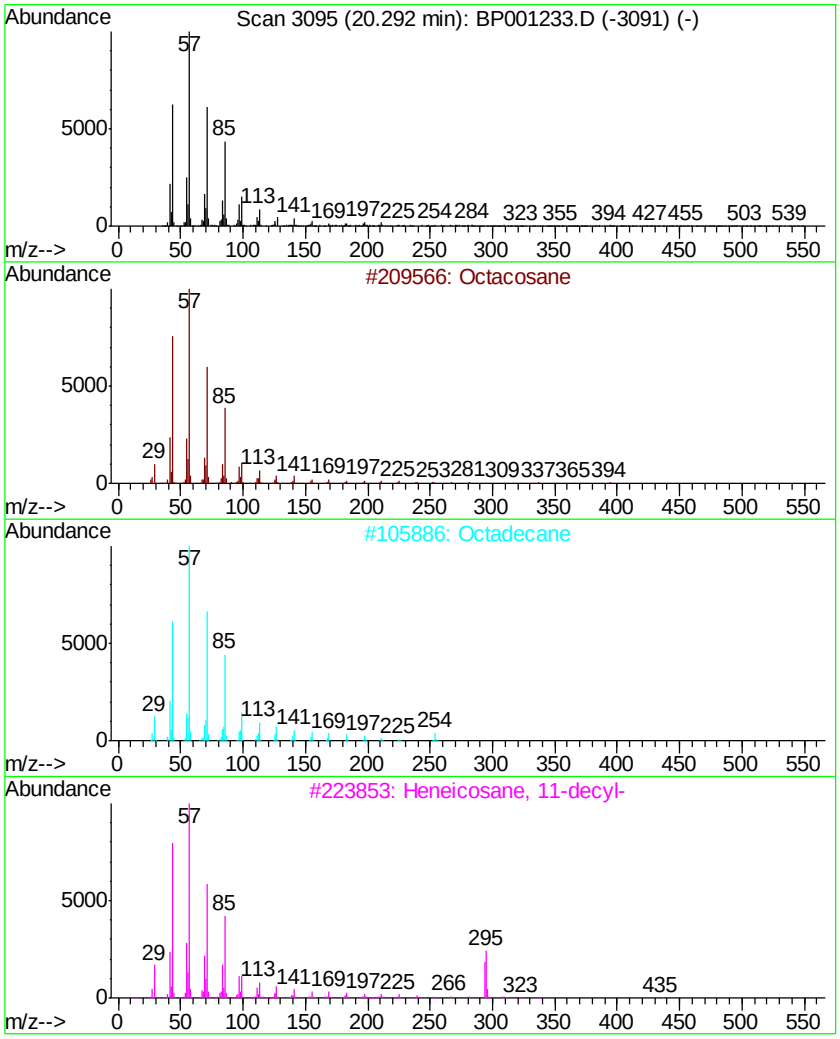
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

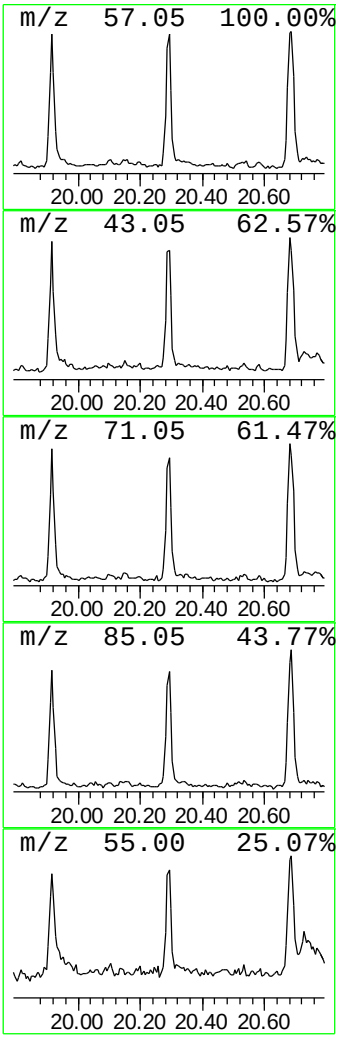
****

Peak Number 8 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.29	3.28 ng	119009	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	94
2	Octadecane	254	C18H38	000593-45-3	93
3	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
4	Hentriacontane	437	C31H64	000630-04-6	91
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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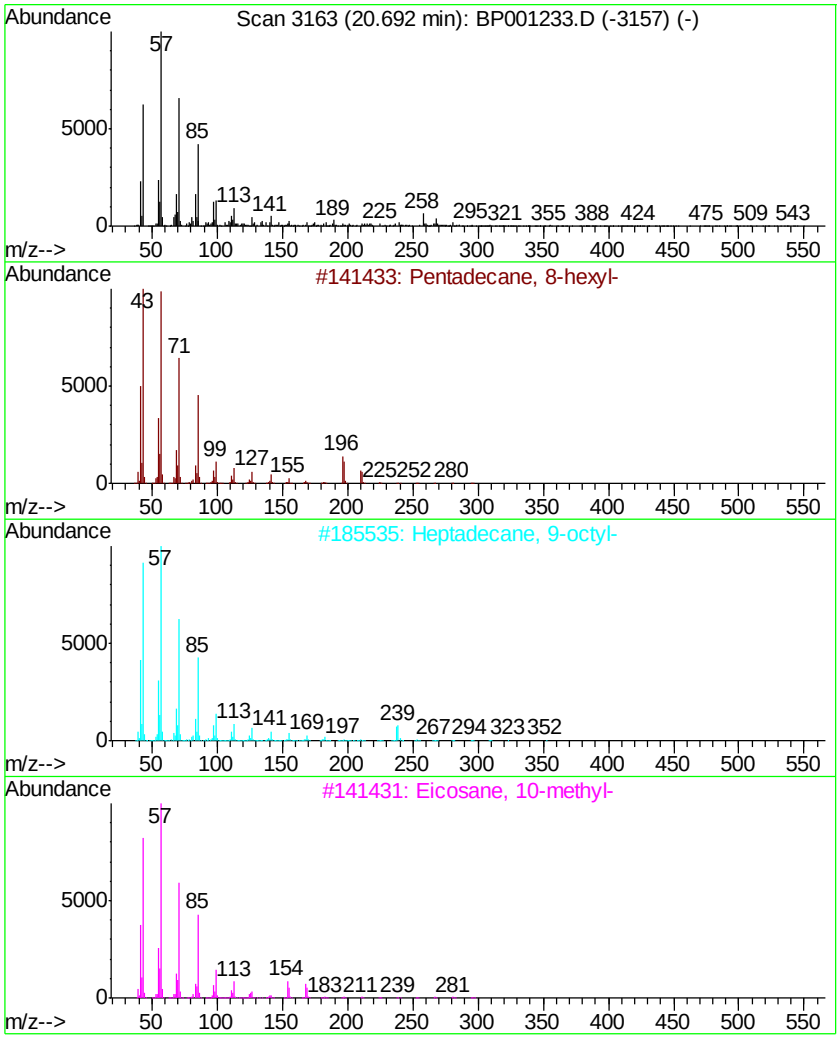
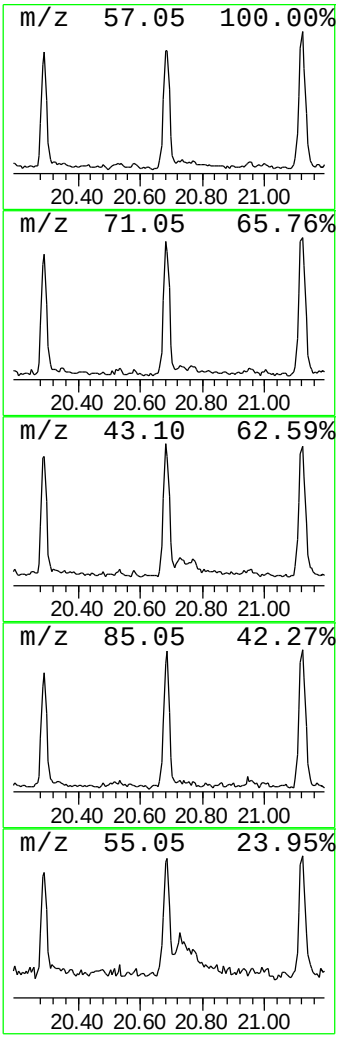
Peak Number 9 Pentadecane, 8-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	4.26 ng	154485	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4	Heneicosane	296	C21H44	000629-94-7	92
5	Nonadecane	268	C19H40	000629-92-5	91





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Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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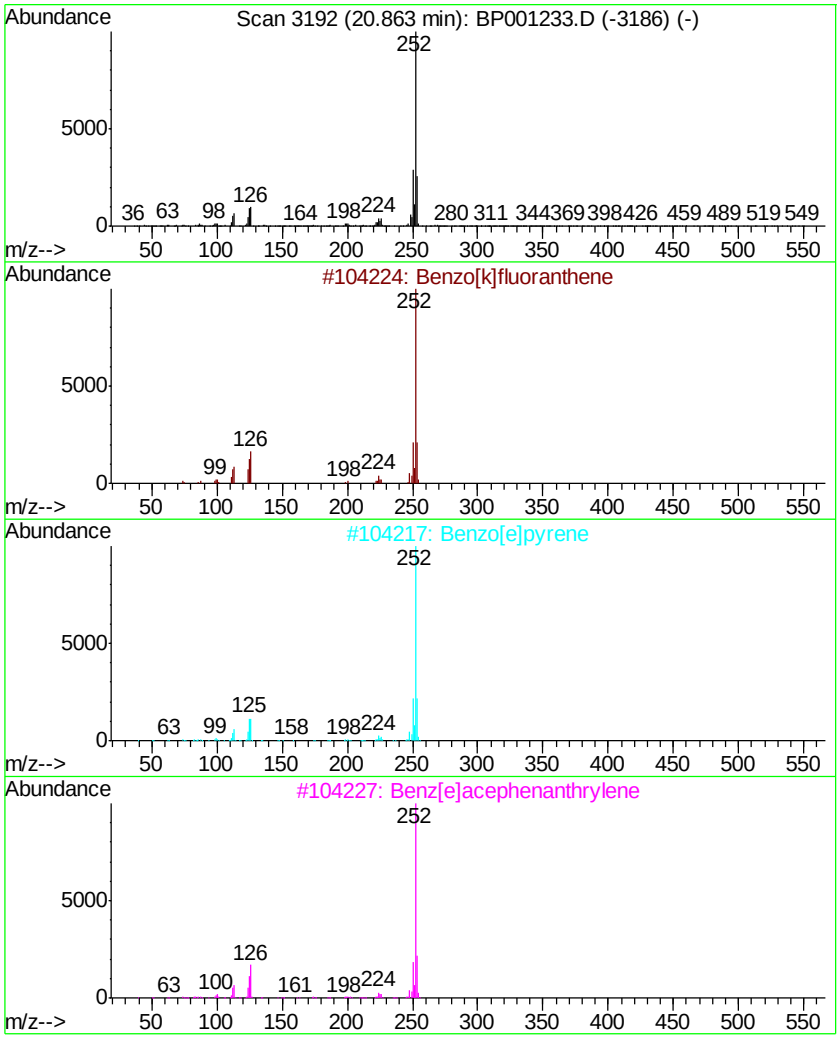
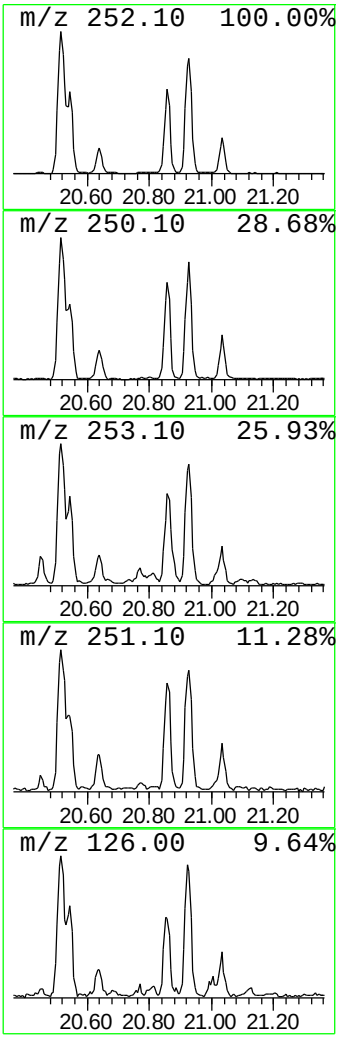
Peak Number 10 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	5.72 ng	207499	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
2	Benzo[e]pyrene	252	C20H12	000192-97-2	94
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
4	Benzo[a]pyrene	252	C20H12	000050-32-8	90
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	86





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Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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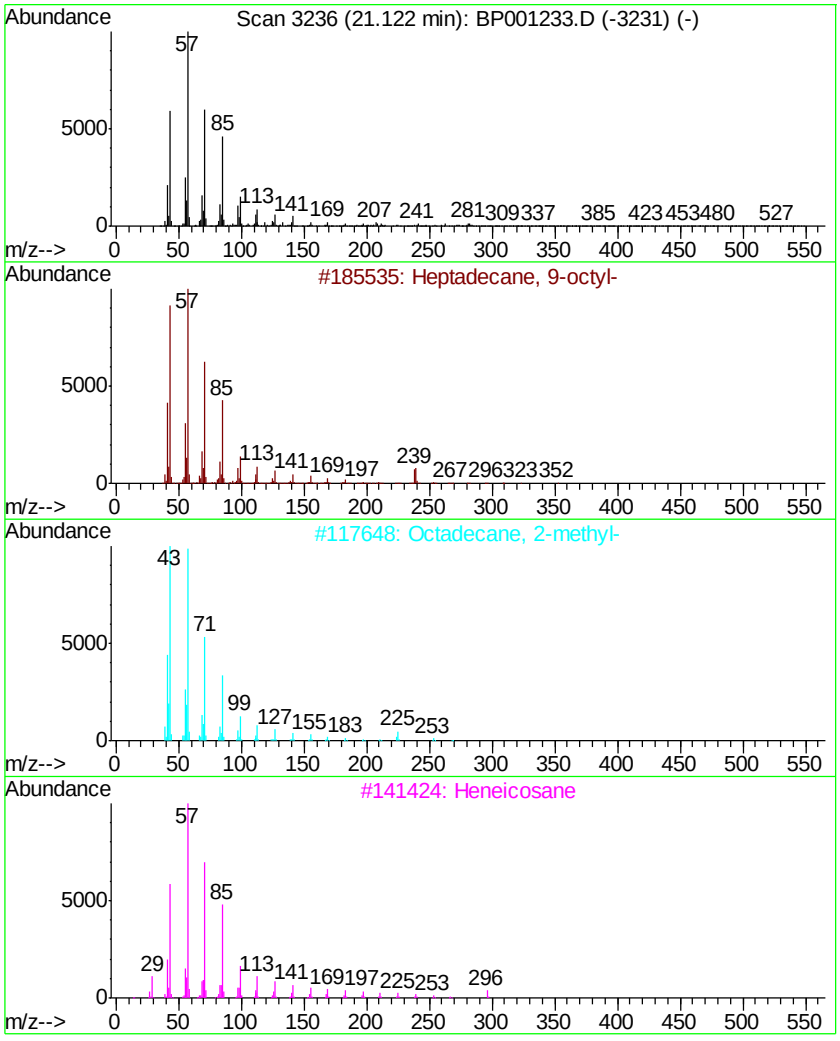
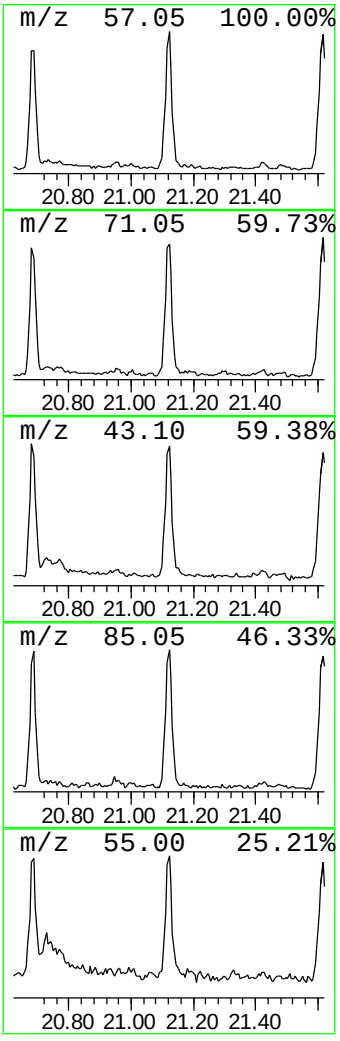
Peak Number 11 Heptadecane, 9-octyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.12	5.20 ng	188498	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2	Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
3	Heneicosane	296	C21H44	000629-94-7	91
4	2-methyloctacosane	408	C29H60	1000376-72-8	90
5	Heptacosane	380	C27H56	000593-49-7	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
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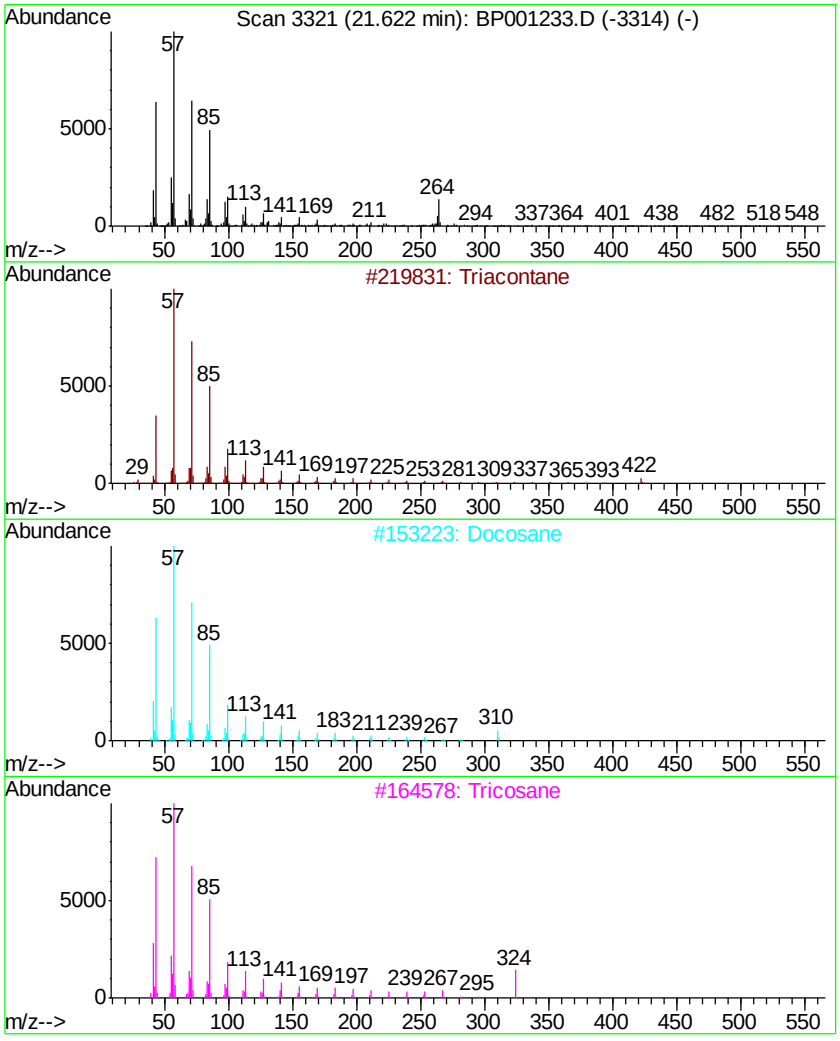
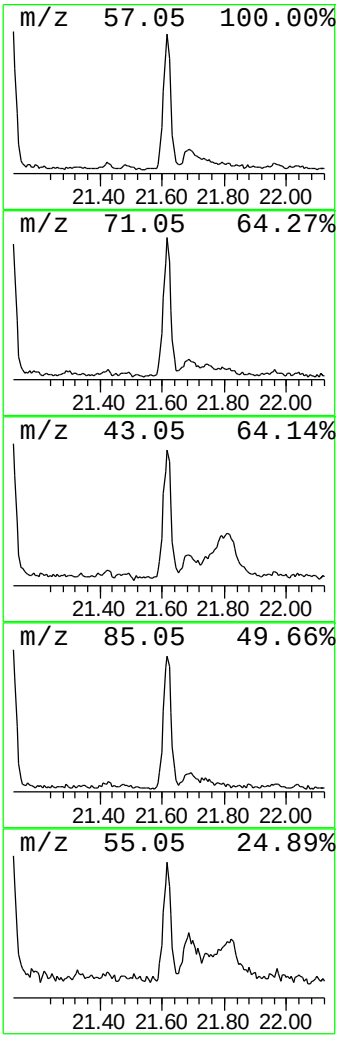
Peak Number 12 Triacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.62	5.74 ng	208176	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	97
2	Docosane	310	C22H46	000629-97-0	95
3	Tricosane	324	C23H48	000638-67-5	95
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5	Pentacosane	352	C25H52	000629-99-2	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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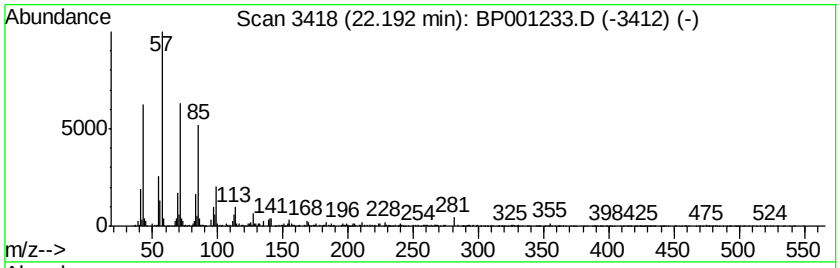
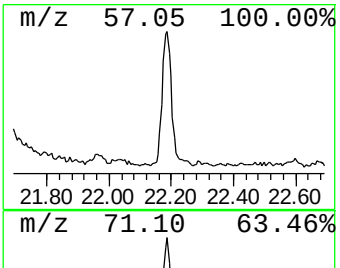
Peak Number 13 Docosane, 11-butyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.19	3.09 ng	112101	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2	Docosane, 9-butyl-	366	C26H54	055282-14-9	90
3	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	90
4	Octacosane	394	C28H58	000630-02-4	90
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87





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Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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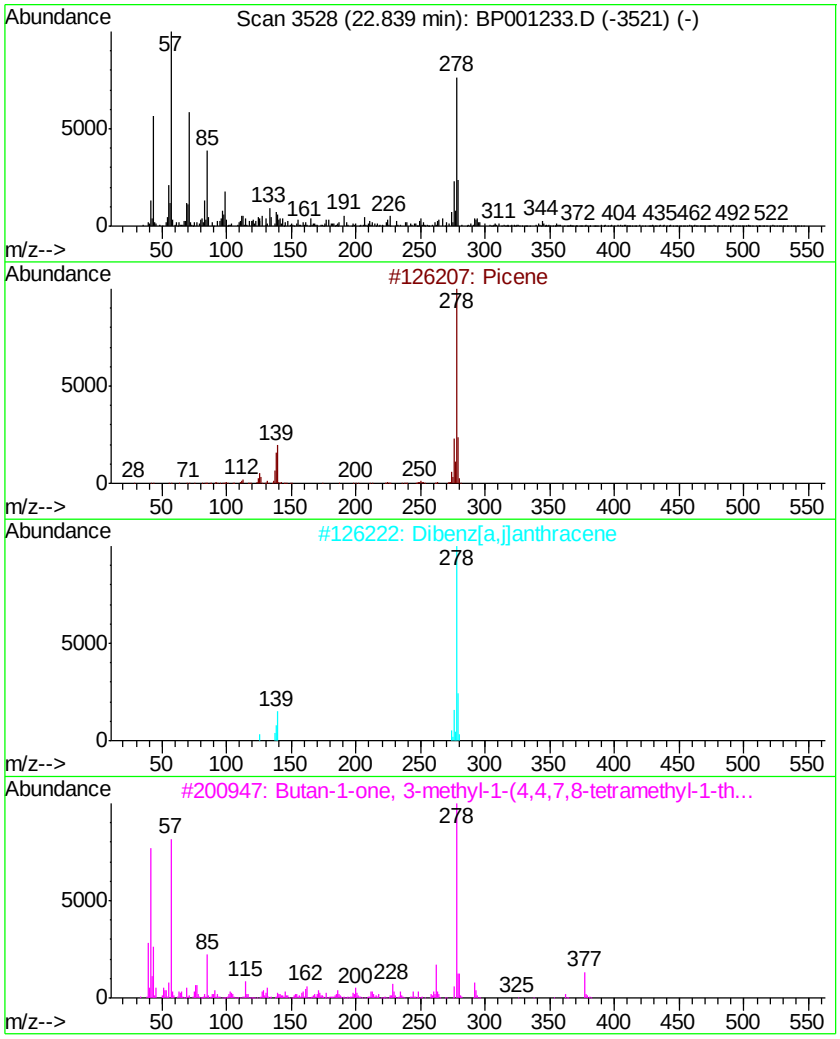
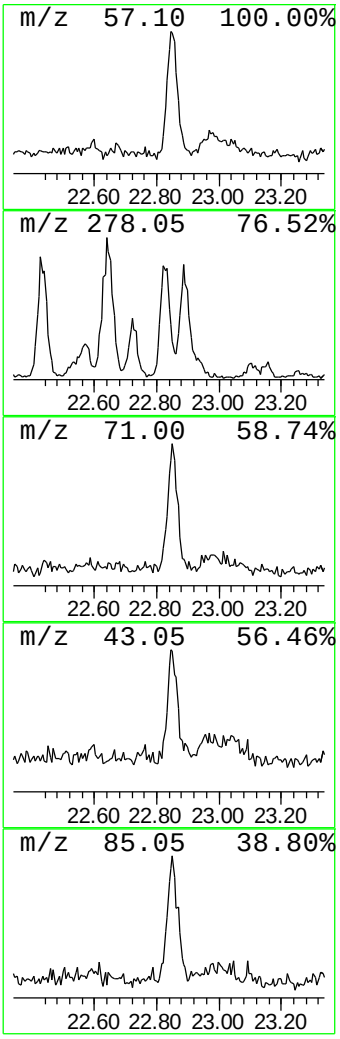
Peak Number 14 Picene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.84	2.75 ng	99500	Perylene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Picene	278	C22H14	000213-46-7	68
2	Dibenz[a,j]anthracene	278	C22H14	000224-41-9	59
3	Butan-1-one, 3-methyl-1-(4,4,7,8...	377	C19H23NOS3	1000310-20-2	47
4	Pentacene	278	C22H14	000135-48-8	46
5	Benzo[b]triphenylene	278	C22H14	000215-58-7	41





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120619\
Data File : BP001233.D
Acq On : 06 Dec 2019 16:54
Operator : JU
Sample : K6119-09
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-2-(0-2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.62	15.7	ng	389549	1	8.30	497338	20.0
unknown7.94	7.94	80.0	ng	1988850	1	8.30	497338	20.0
n-Hexadecanoic acid	16.48	6.0	ng	250058	4	15.59	839956	20.0
Pyrene, 1-methyl-	18.23	2.3	ng	138651	5	19.23	1189990	20.0
Heptafluorobutyri...	19.11	10.4	ng	616561	5	19.23	1189990	20.0
Eicosane, 10-methyl-	19.92	2.1	ng	123608	5	19.23	1189990	20.0
Octacosane	20.29	3.3	ng	119009	6	21.00	724922	20.0
Pentadecane, 8-he...	20.69	4.3	ng	154485	6	21.00	724922	20.0
Benzo[e]pyrene	20.86	5.7	ng	207499	6	21.00	724922	20.0
Heptadecane, 9-oc...	21.12	5.2	ng	188498	6	21.00	724922	20.0
Triacontane	21.62	5.7	ng	208176	6	21.00	724922	20.0
Docosane, 11-butyl-	22.19	3.1	ng	112101	6	21.00	724922	20.0
Picene	22.84	2.8	ng	99500	6	21.00	724922	20.0