

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051017\
 Data File : BF095074.D
 Acq On : 11 May 2017 00:54
 Operator : SJ/MA
 Sample : I2970-01 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-U10-BASE

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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	4.990	516	519	542	rBV	81994	184374	12.92%	0.955%
2	5.654	629	632	652	rBV	400799	921285	64.58%	4.772%
3	5.907	671	675	684	rVB2	30659	49088	3.44%	0.254%
4	7.254	900	904	920	rVV	386763	948809	66.51%	4.914%
5	7.366	920	923	933	rVV	710526	1194650	83.74%	6.188%
6	7.442	933	936	941	rVV	33908	49499	3.47%	0.256%
7	7.701	976	980	990	rVB	467829	556027	38.97%	2.880%
8	7.790	991	995	1001	rVB4	18804	36068	2.53%	0.187%
9	7.937	1016	1020	1033	rBV	698378	911570	63.90%	4.721%
10	8.601	1129	1133	1151	rBV	402494	630008	44.16%	3.263%
11	9.725	1315	1324	1328	rBV	598998	737888	51.72%	3.822%
12	10.519	1453	1459	1464	rBV	33189	40776	2.86%	0.211%
13	10.795	1502	1506	1513	rVB2	26903	31940	2.24%	0.165%
14	10.889	1518	1522	1531	rBV	108460	158672	11.12%	0.822%
15	11.042	1543	1548	1556	rBV	77392	114827	8.05%	0.595%
16	11.448	1612	1617	1621	rVV2	27603	51024	3.58%	0.264%
17	11.495	1621	1625	1640	rVB	1154664	1426628	100.00%	7.389%
18	11.719	1659	1663	1671	rVB	35908	45818	3.21%	0.237%
19	11.913	1692	1696	1698	rBV3	24063	36674	2.57%	0.190%
20	12.030	1712	1716	1720	rBV	46446	64010	4.49%	0.332%
21	12.072	1720	1723	1731	rVB2	45042	70842	4.97%	0.367%
22	12.195	1739	1744	1749	rBV	129153	197394	13.84%	1.022%
23	12.236	1749	1751	1755	rVB3	45999	54836	3.84%	0.284%
24	12.330	1764	1767	1778	rVV2	54084	87979	6.17%	0.456%
25	12.554	1800	1805	1815	rBV2	643199	932064	65.33%	4.828%
26	12.766	1835	1841	1848	rBV3	13327	31749	2.23%	0.164%
27	12.907	1860	1865	1867	rBV	37249	49416	3.46%	0.256%
28	13.077	1888	1894	1896	rVV4	27796	44085	3.09%	0.228%
29	13.101	1896	1898	1904	rVV3	41889	56870	3.99%	0.295%
30	13.395	1942	1948	1954	rBV	67310	92836	6.51%	0.481%
31	13.454	1954	1958	1966	rVB2	60753	90459	6.34%	0.469%
32	13.730	2000	2005	2007	rBV	29327	38999	2.73%	0.202%
33	13.754	2007	2009	2014	rVB	38548	46533	3.26%	0.241%
34	13.854	2021	2026	2035	rBV	431730	693825	48.63%	3.594%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.924	2035	2038	2044	rVB7	22741	34320	2.41%	0.178%
36	14.166	2076	2079	2081	rBV	61598	73689	5.17%	0.382%
37	14.189	2081	2083	2088	rVV	100829	129713	9.09%	0.672%
38	14.489	2127	2134	2140	rBV10	13741	35469	2.49%	0.184%
39	14.577	2145	2149	2154	rVV4	33047	60322	4.23%	0.312%
40	14.683	2164	2167	2172	rVV4	27863	51177	3.59%	0.265%
41	14.754	2177	2179	2183	rVV3	24825	37891	2.66%	0.196%
42	14.901	2200	2204	2208	rBV	56070	63347	4.44%	0.328%
43	14.954	2208	2213	2217	rBV	663616	817457	57.30%	4.234%
44	14.995	2217	2220	2227	rVB	156288	214894	15.06%	1.113%
45	15.083	2231	2235	2242	rVB	47406	61349	4.30%	0.318%
46	15.177	2245	2251	2255	rBV8	21453	38704	2.71%	0.200%
47	15.483	2299	2303	2307	rBV6	23302	34095	2.39%	0.177%
48	15.560	2313	2316	2321	rVB	58457	70931	4.97%	0.367%
49	15.742	2342	2347	2350	rBV	68264	77291	5.42%	0.400%
50	15.783	2350	2354	2361	rVB	94912	123979	8.69%	0.642%
51	15.895	2369	2373	2375	rVV	75704	92032	6.45%	0.477%
52	15.924	2375	2378	2383	rVV2	122395	200549	14.06%	1.039%
53	15.965	2383	2385	2388	rVB	34367	31466	2.21%	0.163%
54	16.154	2413	2417	2420	rBV	55913	68769	4.82%	0.356%
55	16.183	2420	2422	2429	rVV3	50533	72909	5.11%	0.378%
56	16.242	2429	2432	2437	rVB5	23875	35157	2.46%	0.182%
57	16.436	2462	2465	2469	rBV2	30219	36140	2.53%	0.187%
58	16.536	2478	2482	2486	rBV	101732	129932	9.11%	0.673%
59	16.612	2493	2495	2498	rVB2	34631	33413	2.34%	0.173%
60	16.660	2499	2503	2507	rBV2	46025	62048	4.35%	0.321%
61	16.701	2507	2510	2513	rBV	52808	59189	4.15%	0.307%
62	16.742	2513	2517	2522	rBV	250987	305424	21.41%	1.582%
63	16.836	2529	2533	2536	rBV2	28561	37648	2.64%	0.195%
64	17.012	2560	2563	2564	rVV3	41315	39466	2.77%	0.204%
65	17.042	2564	2568	2576	rVB	314042	384914	26.98%	1.994%
66	17.159	2585	2588	2592	rVB2	31873	39768	2.79%	0.206%
67	17.201	2592	2595	2598	rBV	49143	50781	3.56%	0.263%
68	17.277	2603	2608	2616	rVB	893054	1040341	72.92%	5.388%
69	17.359	2619	2622	2626	rBV2	22716	38271	2.68%	0.198%
70	17.477	2638	2642	2644	rBV4	26126	36914	2.59%	0.191%
71	17.642	2666	2670	2673	rBV2	64487	101678	7.13%	0.527%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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72	17.671	2673	2675	2681	rVB	54583	55720	3.91%	0.289%
73	17.754	2686	2689	2693	rVB	42156	48882	3.43%	0.253%
74	18.112	2747	2750	2753	rVB2	43574	38266	2.68%	0.198%
75	18.183	2759	2762	2767	rBV	42091	47100	3.30%	0.244%
76	18.306	2780	2783	2785	rBV2	35116	39234	2.75%	0.203%
77	18.336	2785	2788	2790	rVB2	35953	43214	3.03%	0.224%
78	18.536	2818	2822	2825	rBV	57593	70326	4.93%	0.364%
79	18.624	2832	2837	2841	rBV2	641995	868029	60.84%	4.496%
80	18.659	2841	2843	2849	rVB	143820	182501	12.79%	0.945%
81	18.753	2856	2859	2862	rVB4	39908	55111	3.86%	0.285%
82	19.136	2918	2924	2927	rVB2	58526	78855	5.53%	0.408%
83	19.171	2928	2930	2934	rVB	34260	35300	2.47%	0.183%
84	19.318	2953	2955	2958	rVB2	45040	42905	3.01%	0.222%
85	19.548	2992	2994	3002	rVB8	26227	50570	3.54%	0.262%
86	19.736	3022	3026	3030	rBV2	48504	75739	5.31%	0.392%
87	19.895	3048	3053	3056	rBV	222396	330386	23.16%	1.711%
88	19.918	3056	3057	3063	rVB	119302	96645	6.77%	0.501%
89	20.000	3069	3071	3075	rBV	59916	66648	4.67%	0.345%
90	20.036	3075	3077	3081	rVB3	41722	38804	2.72%	0.201%
91	20.183	3098	3102	3108	rBV	159285	201071	14.09%	1.041%
92	20.242	3108	3112	3116	rBV	177982	190004	13.32%	0.984%
93	20.295	3117	3121	3125	rBV	533678	590805	41.41%	3.060%
94	20.383	3134	3136	3142	rVB5	38875	51369	3.60%	0.266%
95	20.759	3197	3200	3204	rVB2	46824	49763	3.49%	0.258%
96	21.183	3269	3272	3277	rVB4	23923	38042	2.67%	0.197%
97	21.595	3337	3342	3350	rVB3	106736	201773	14.14%	1.045%
98	21.671	3350	3355	3361	rVB8	21050	41109	2.88%	0.213%
99	21.742	3363	3367	3371	rVB2	23793	35929	2.52%	0.186%
100	21.971	3401	3406	3415	rVB2	75618	144118	10.10%	0.746%

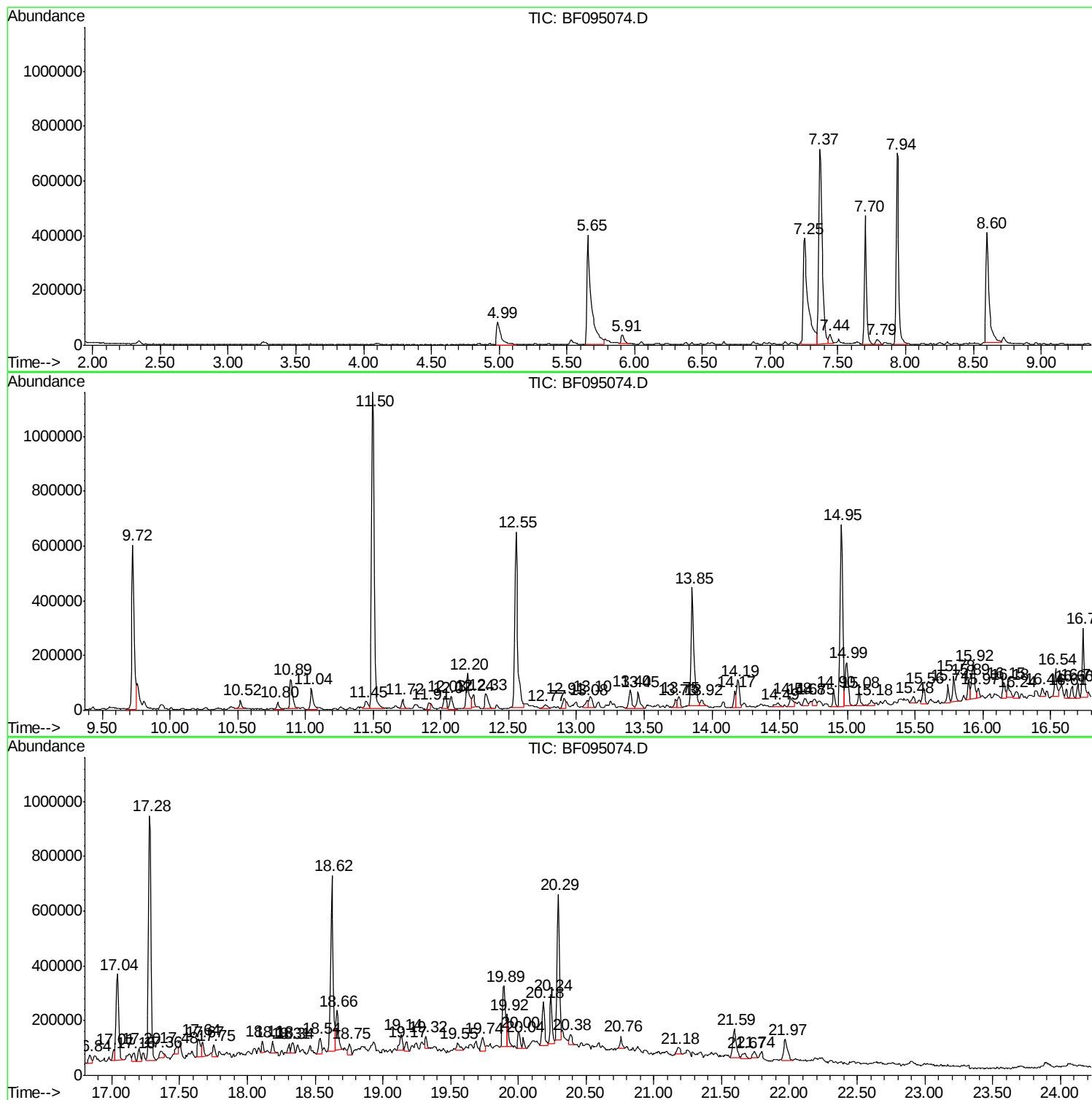
Sum of corrected areas: 19307177

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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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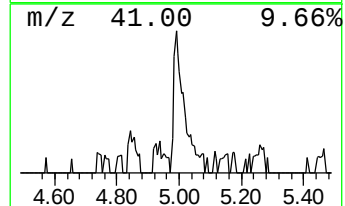
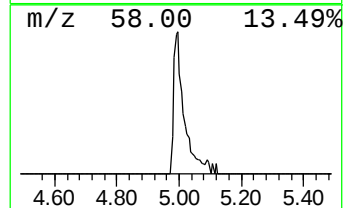
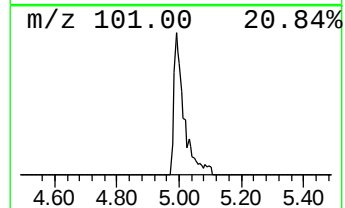
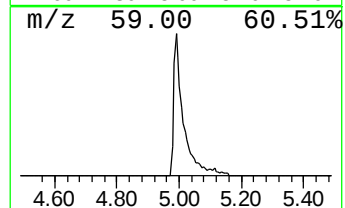
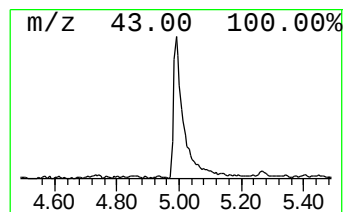
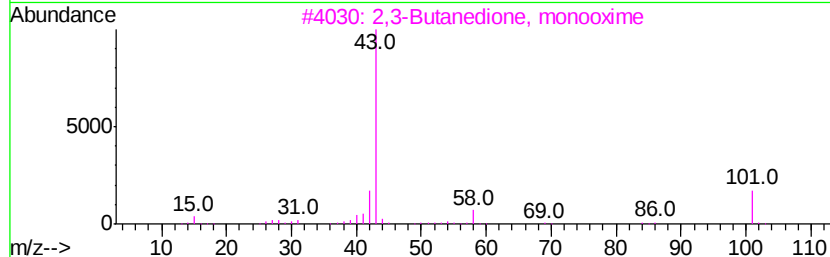
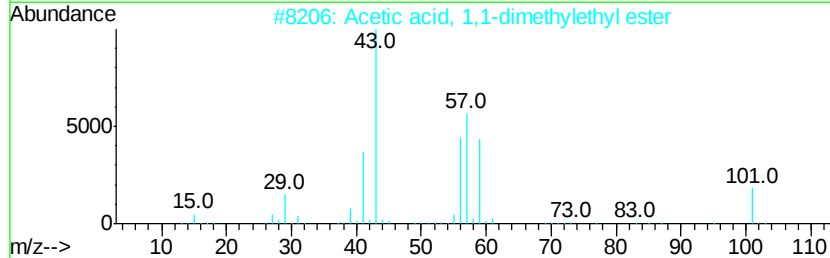
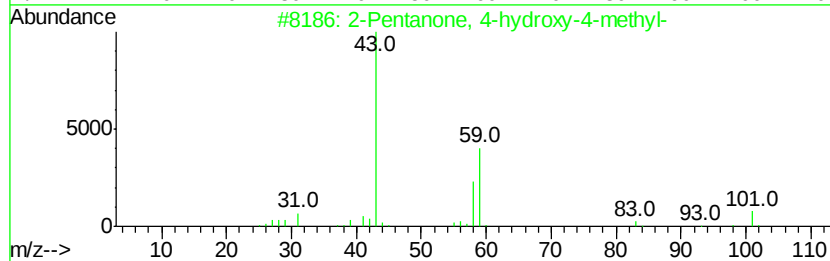
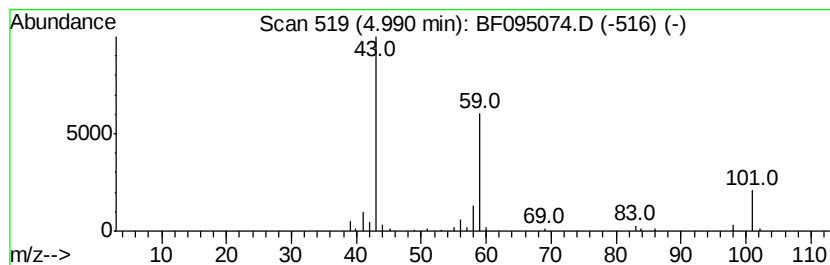
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	6.63 ng	184374	1,4-Dichlorobenzene-d4	7.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	38
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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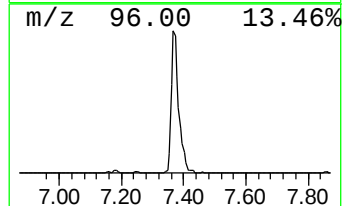
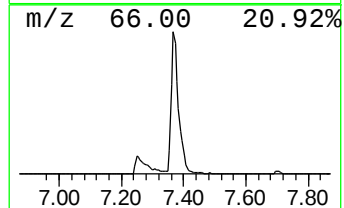
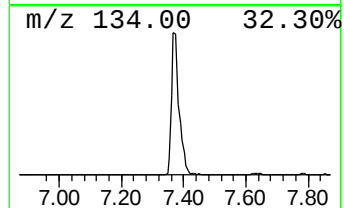
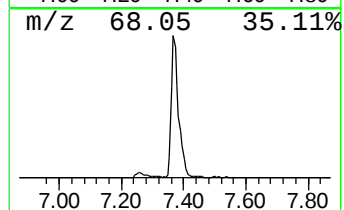
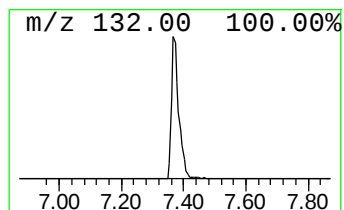
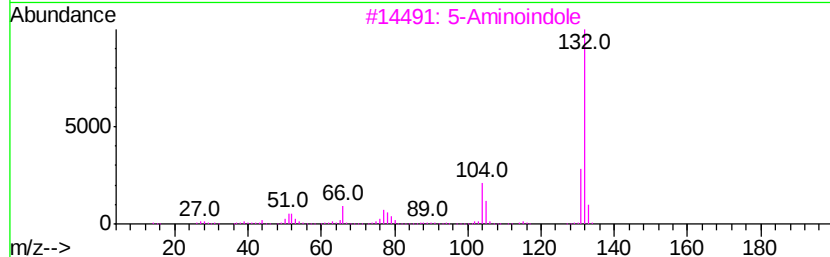
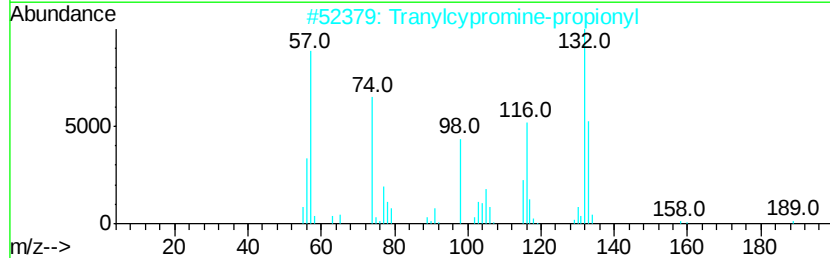
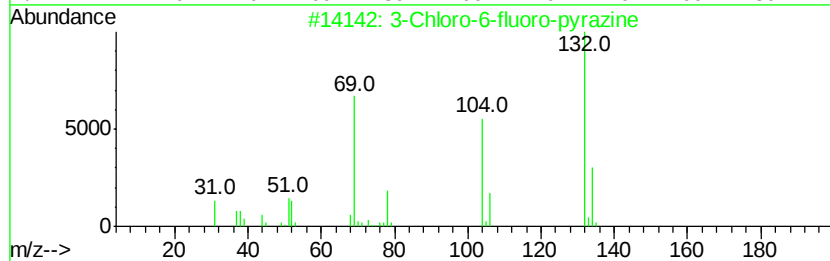
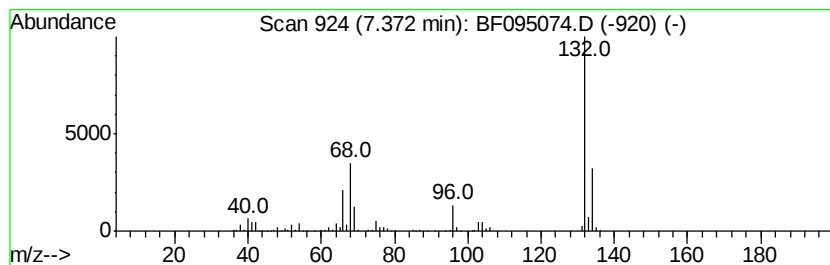
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.37 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	42.97 ng	1194650	1,4-Dichlorobenzene-d4	7.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5		Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	12



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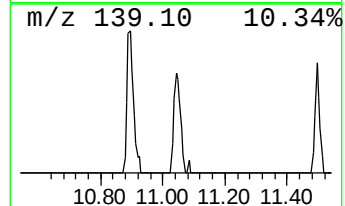
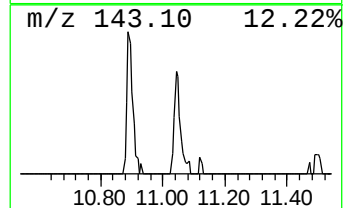
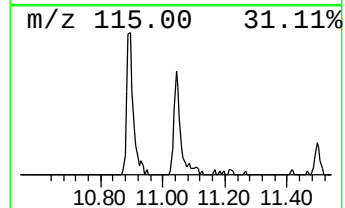
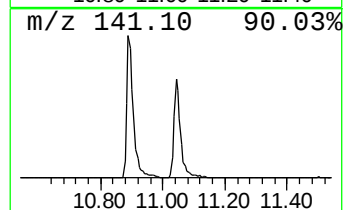
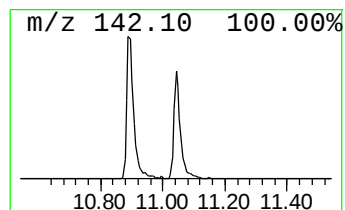
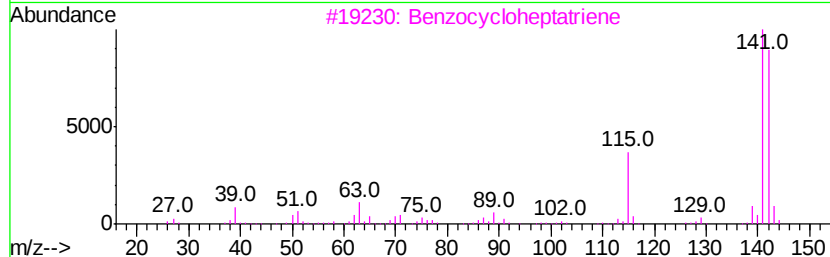
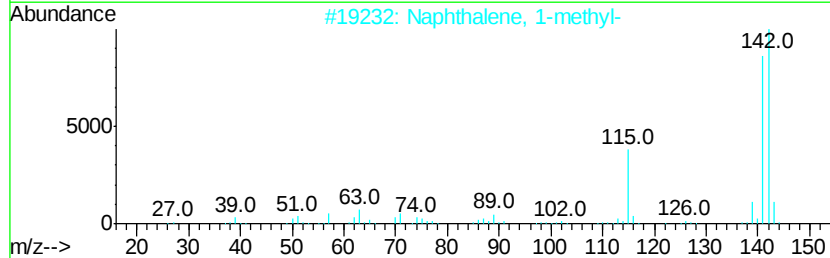
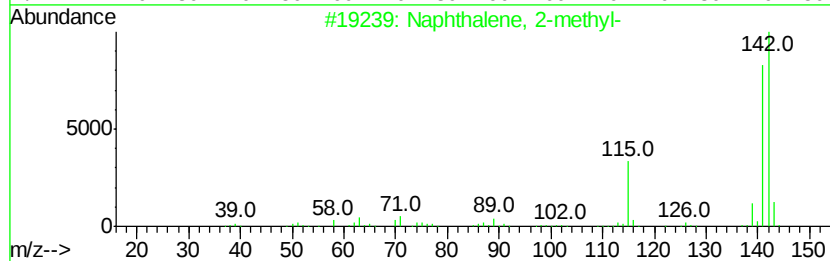
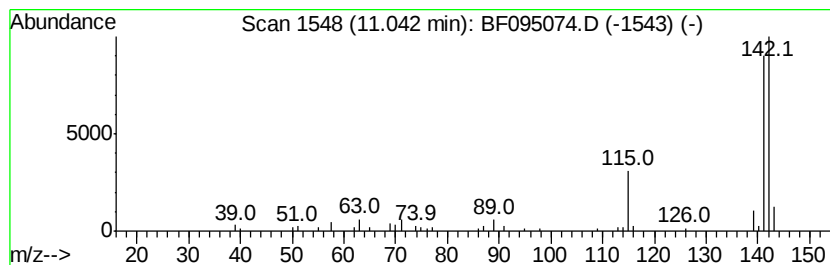
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.04	3.11 ng	114827	Naphthalene-d8	9.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5			3,4-Difluorobenzaldehyde	142	C7H4F2O	034036-07-2	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
Sample : I2970-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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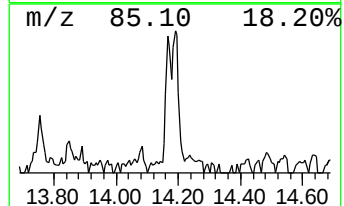
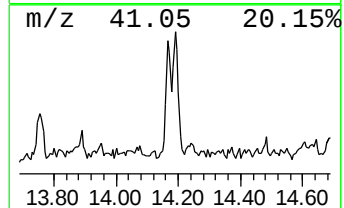
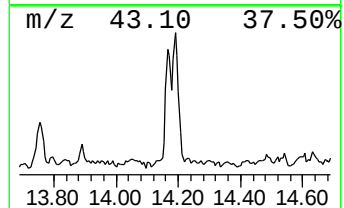
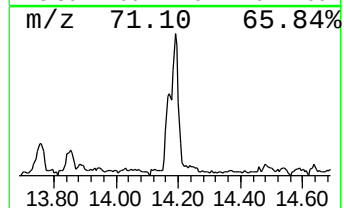
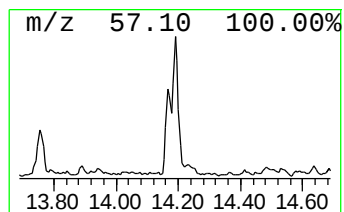
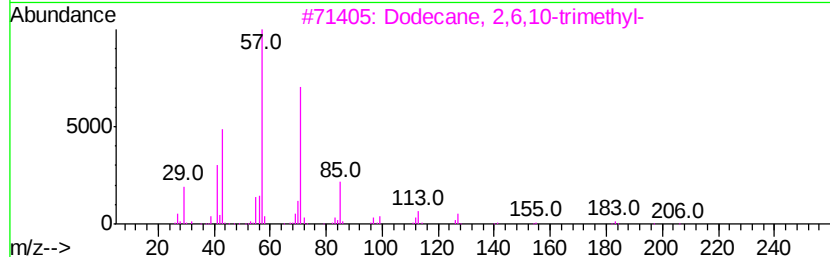
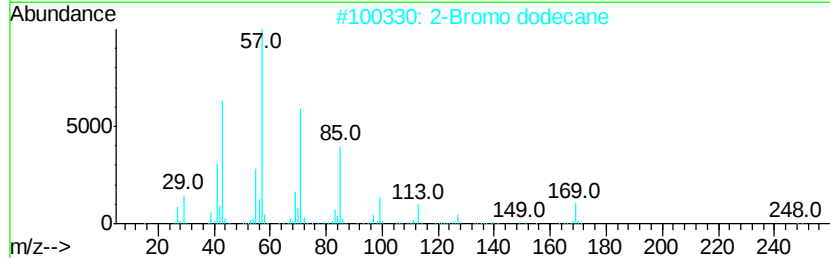
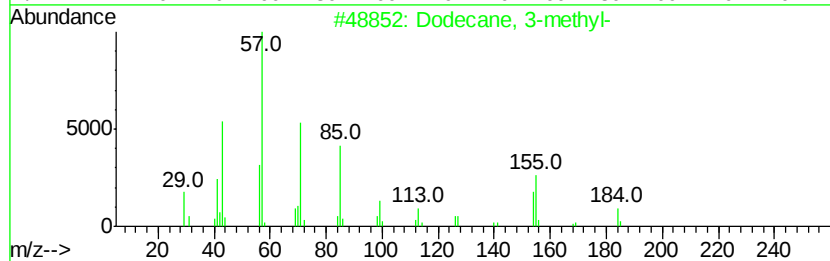
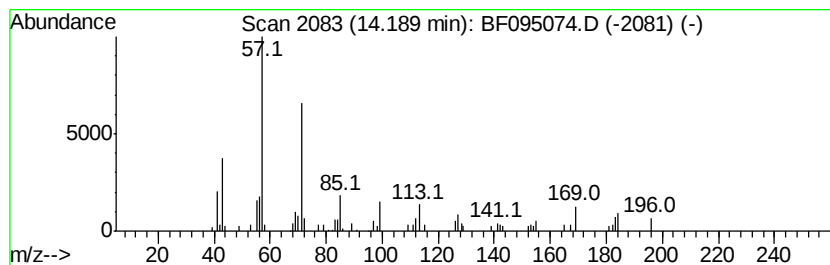
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Dodecane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	3.17 ng	129713	Phenanthrene-d10	14.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 3-methyl-	184	C13H28	017312-57-1	58	
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	58	
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	53	
4	Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	50	
5	Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	50	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
Sample : I2970-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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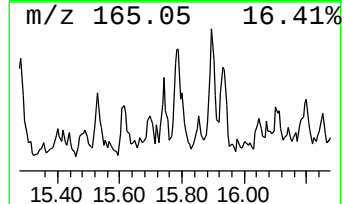
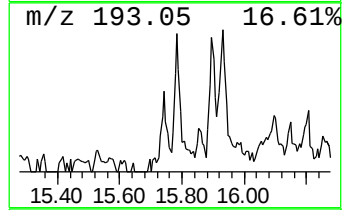
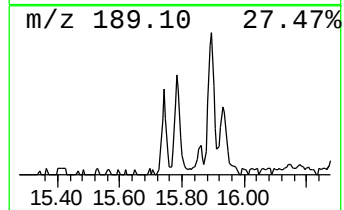
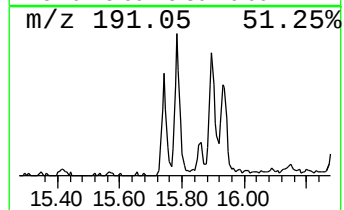
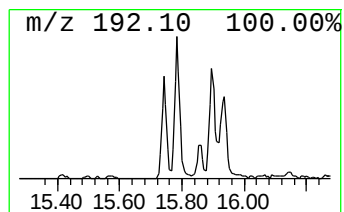
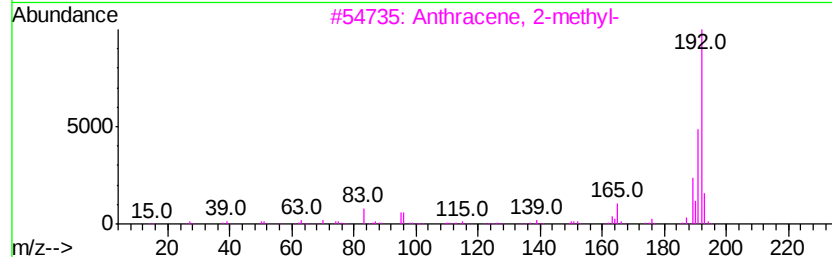
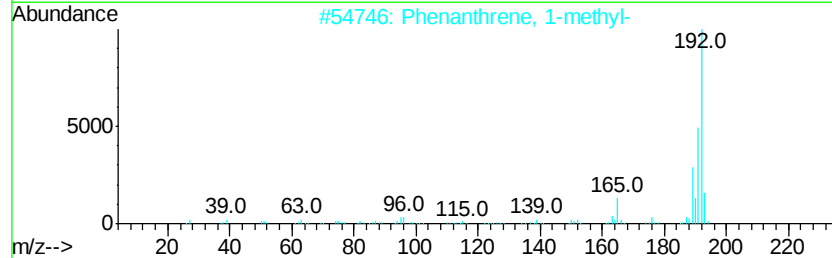
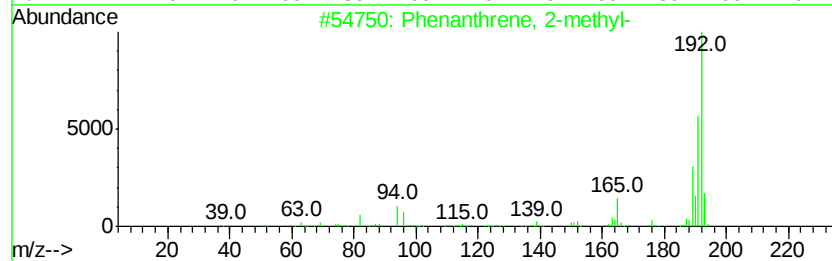
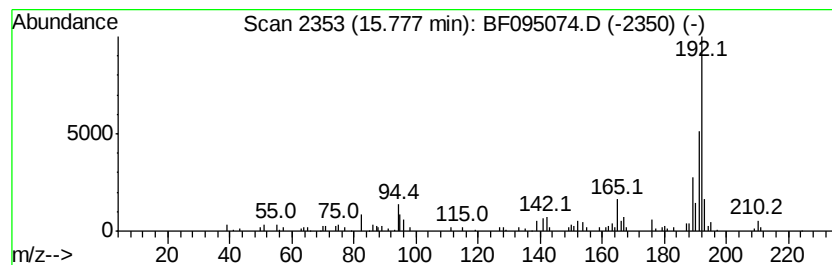
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.78	3.03 ng	123979	Phenanthrene-d10	14.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	95
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
Sample : I2970-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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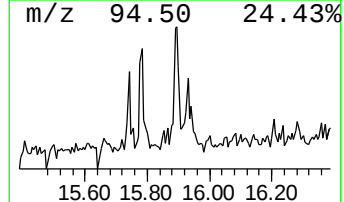
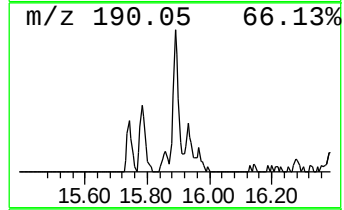
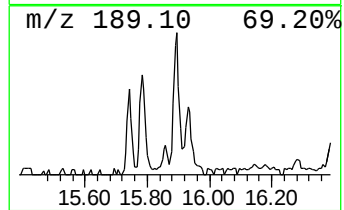
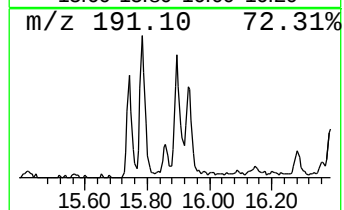
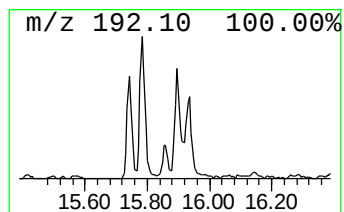
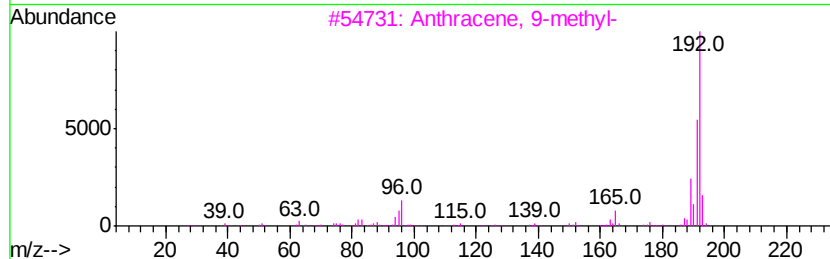
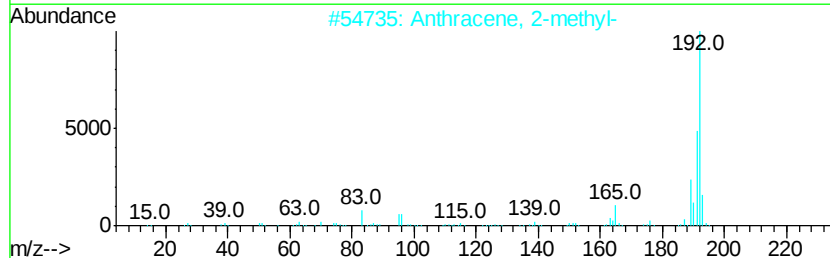
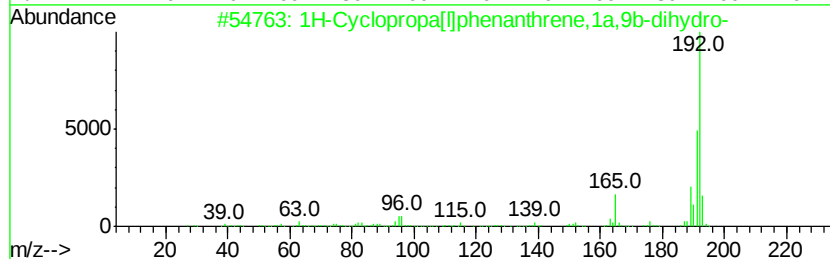
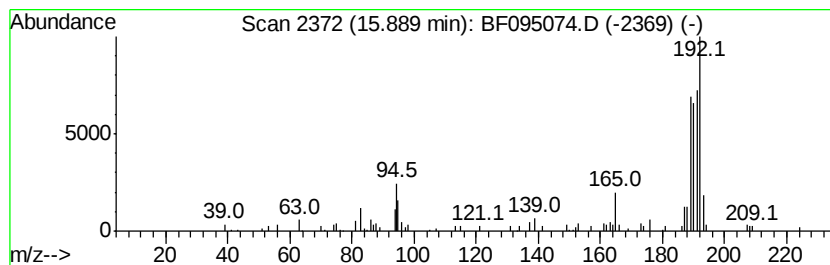
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Cyclopropa[l]phenanthren... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.89	2.25 ng	92032	Phenanthrene-d10	14.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	76
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	70
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	70
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	68
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
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Misc :
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Instrument :
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ClientSampleId :
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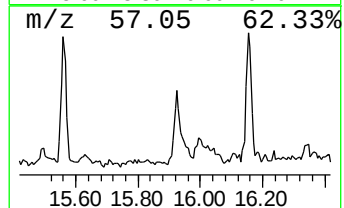
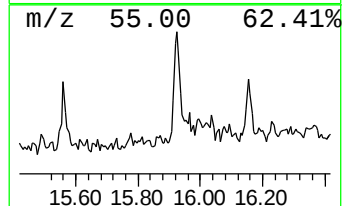
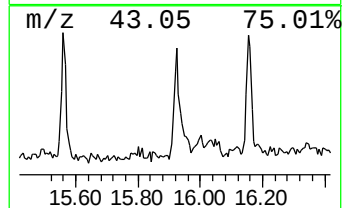
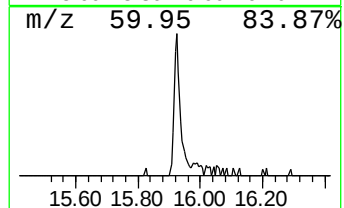
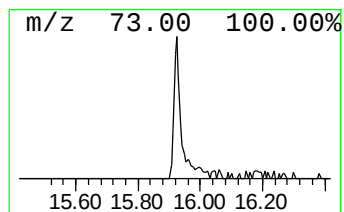
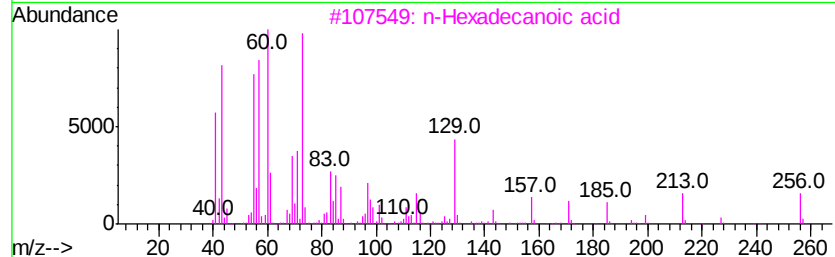
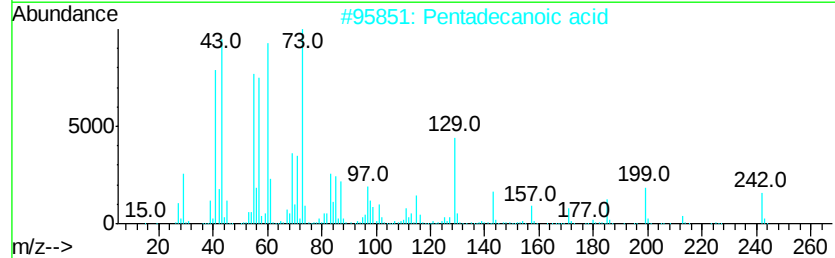
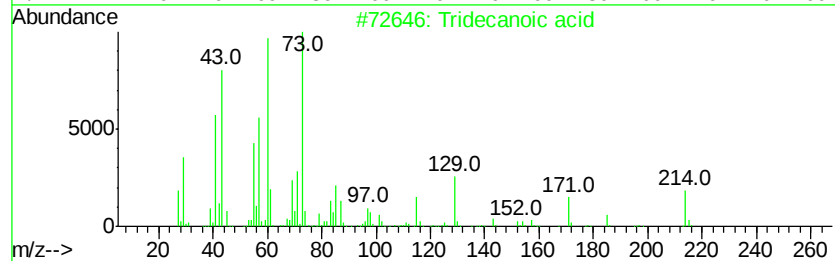
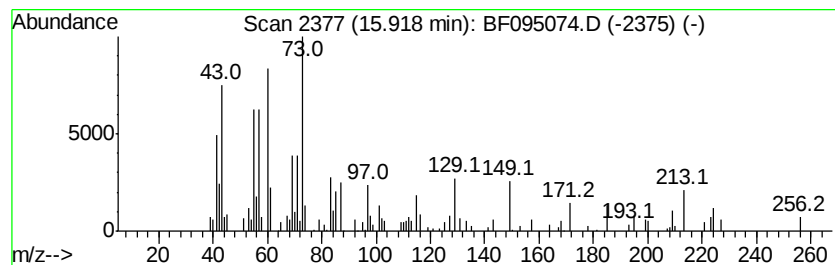
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	4.91 ng	200549	Phenanthrene-d10	14.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	60
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	49
5	Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
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Misc :
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Instrument :
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ClientSampled :
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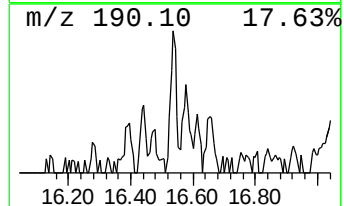
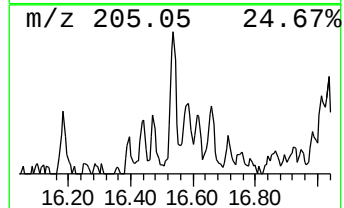
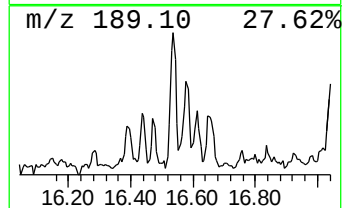
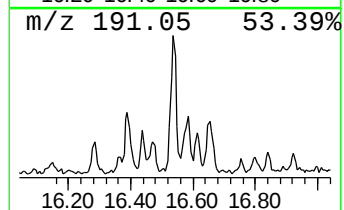
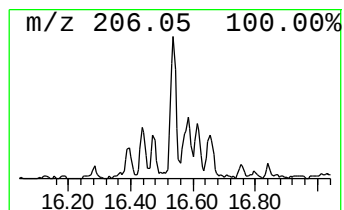
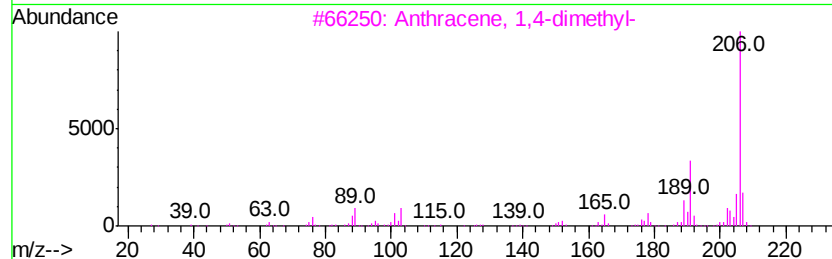
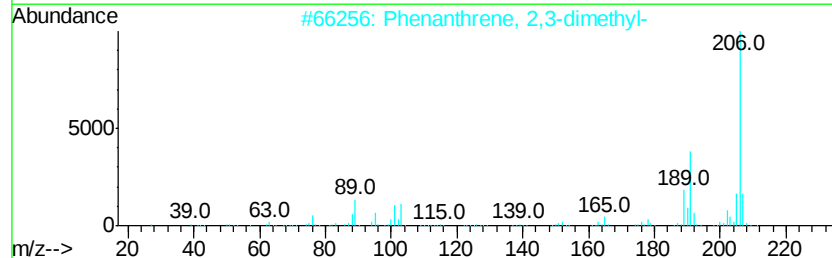
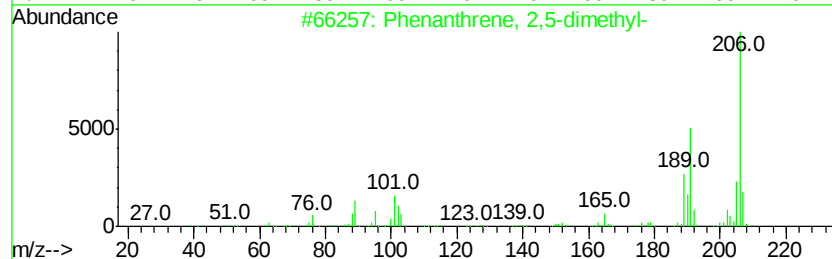
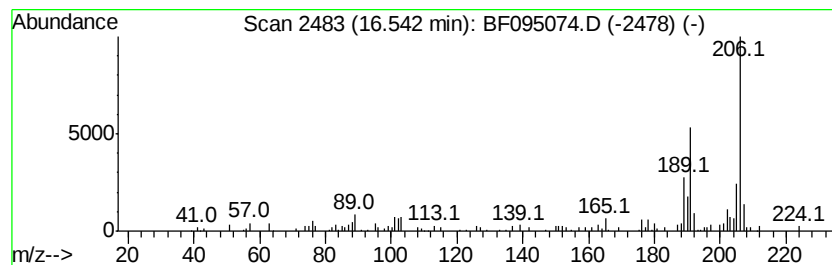
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	3.18 ng	129932	Phenanthrene-d10	14.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94
3		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	92
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
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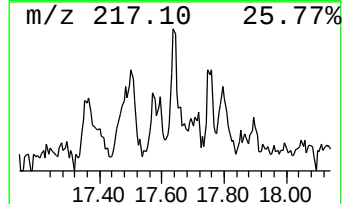
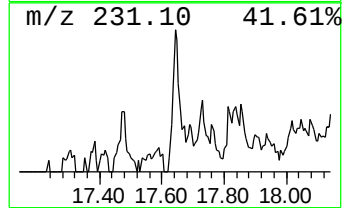
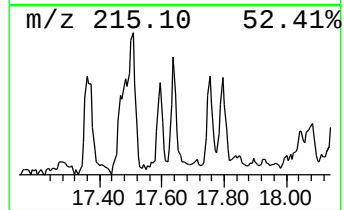
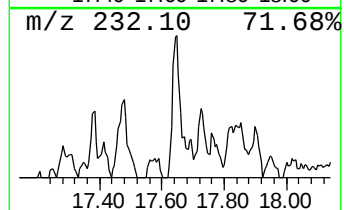
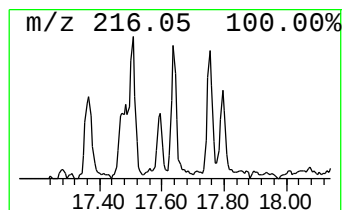
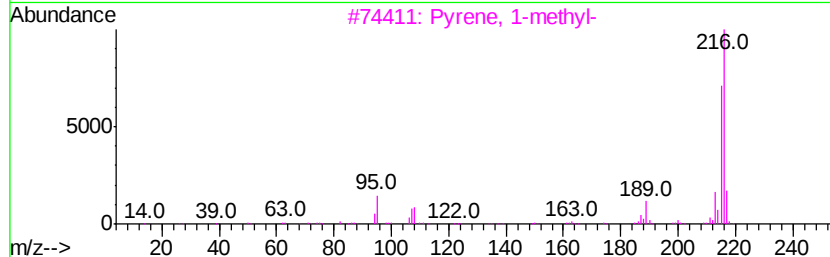
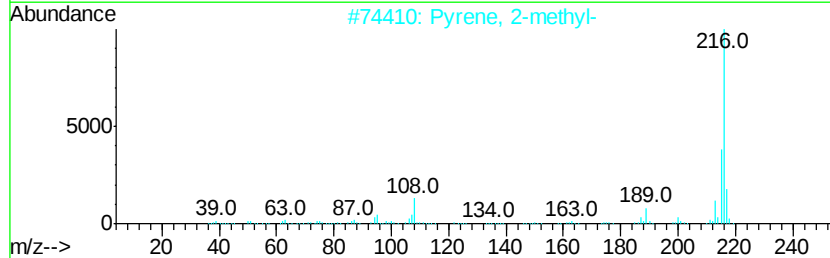
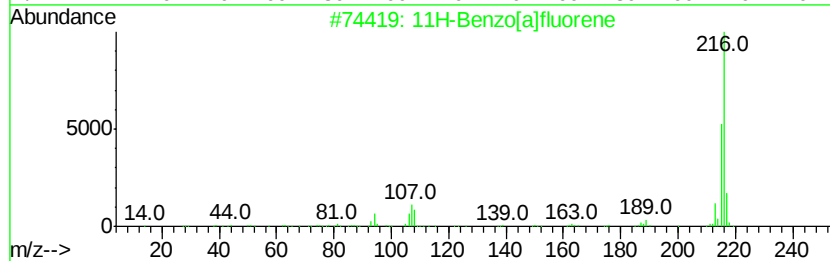
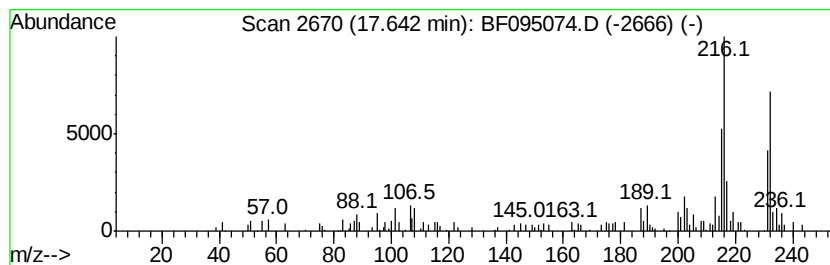
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.64	2.34 ng	101678	Chrysene-d12	18.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	50
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	50
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	50
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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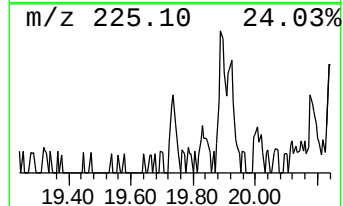
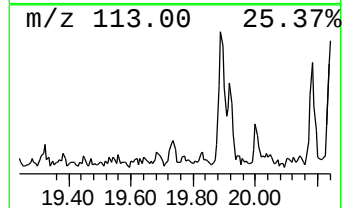
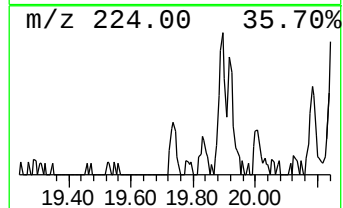
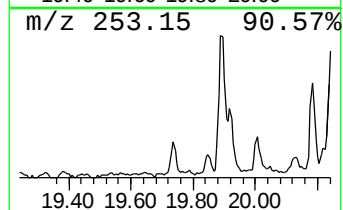
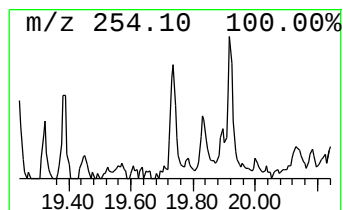
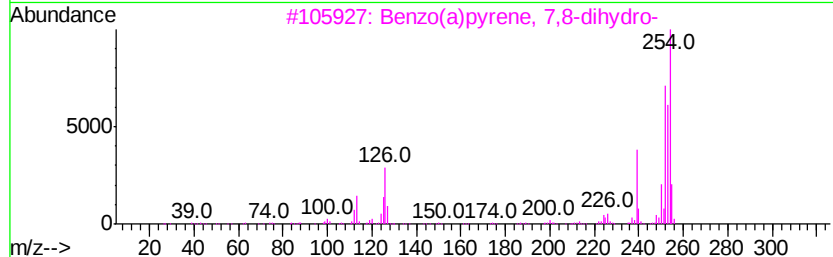
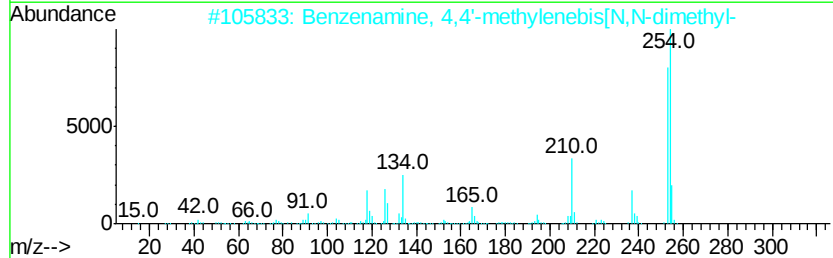
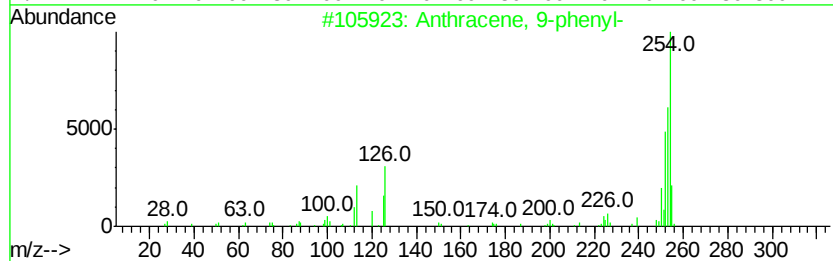
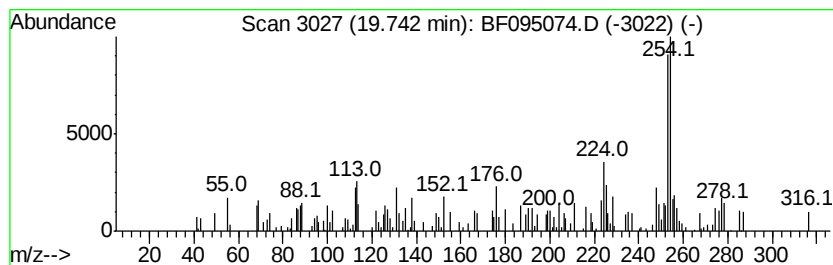
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 9-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.74	2.56 ng	75739	Perylene-d12	20.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-phenyl-	254	C20H14	000602-55-1	70
2		Benzenamine, 4,4'-methylenebis[N...	254	C17H22N2	000101-61-1	64
3		Benzo(a)pyrene, 7,8-dihydro-	254	C20H14	017573-23-8	58
4		Benzenamine, 4-(2-benzothiazolyl...	254	C15H14N2S	010205-56-8	58
5		1,2'-Binaphthalene	254	C20H14	004325-74-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051017\
Data File : BF095074.D
Acq On : 11 May 2017 00:54
Operator : SJ/MA
Sample : I2970-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-U10-BASE

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.99	6.6	ng	184374	1	7.70	556027	20.0
unknown7.37	7.37	43.0	ng	1194650	1	7.70	556027	20.0
Naphthalene, 1-me...	11.04	3.1	ng	114827	2	9.72	737888	20.0
Dodecane, 3-methyl-	14.19	3.2	ng	129713	4	14.95	817457	20.0
Phenanthrene, 2-m...	15.78	3.0	ng	123979	4	14.95	817457	20.0
1H-Cyclopropa[l]p...	15.89	2.3	ng	92032	4	14.95	817457	20.0
Tridecanoic acid	15.92	4.9	ng	200549	4	14.95	817457	20.0
Phenanthrene, 2,5...	16.54	3.2	ng	129932	4	14.95	817457	20.0
11H-Benzo[a]fluorene	17.64	2.3	ng	101678	5	18.62	868029	20.0
Anthracene, 9-phe...	19.74	2.6	ng	75739	6	20.29	590805	20.0