

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010224\
 Data File : BG060211.D
 Acq On : 3 Jan 2024 3:16
 Operator : MA/JU
 Sample : 05899-24
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GHL-SS-19-2-4

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_G\Methods\8270-BG122923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG060211.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	10	rVB	428338	443259	2.90%	0.487%
2	5.508	405	412	429	rBV	3660831	5827700	38.13%	6.403%
3	7.100	675	683	699	rBV	3568081	6196262	40.54%	6.808%
4	7.935	818	825	834	rBV	827055	1441756	9.43%	1.584%
5	9.098	1015	1023	1034	rBV	2029706	3651302	23.89%	4.012%
6	10.743	1294	1303	1310	rBV	1251810	2277443	14.90%	2.502%
7	13.199	1713	1721	1738	rBV	6454214	10306247	67.43%	11.323%
8	14.574	1948	1955	1962	rBV	2516405	3811995	24.94%	4.188%
9	16.060	2199	2208	2221	rBV	3174072	4916442	32.17%	5.402%
10	17.324	2416	2423	2428	rBV	3306024	5040228	32.98%	5.538%
11	17.365	2428	2430	2435	rVB	220575	264763	1.73%	0.291%
12	17.453	2439	2445	2451	rBV	1793189	2761610	18.07%	3.034%
13	17.600	2464	2470	2475	rBV5	142496	209699	1.37%	0.230%
14	17.723	2483	2491	2498	rBV	244517	428862	2.81%	0.471%
15	18.211	2570	2574	2580	rBV2	287090	477819	3.13%	0.525%
16	18.270	2580	2584	2589	rVV	168304	275955	1.81%	0.303%
17	18.328	2590	2594	2600	rVB	174479	250569	1.64%	0.275%
18	19.374	2766	2772	2782	rBV	1376985	2376222	15.55%	2.611%
19	19.492	2788	2792	2797	rVV2	136711	175494	1.15%	0.193%
20	19.562	2799	2804	2811	rVV	595713	880884	5.76%	0.968%
21	19.738	2824	2834	2842	rBV	969235	1549188	10.14%	1.702%
22	19.944	2859	2869	2882	rBV	11282629	15284131	100.00%	16.792%
23	20.067	2885	2890	2895	rVV6	76223	183663	1.20%	0.202%
24	20.990	3043	3047	3053	rVB	141632	202955	1.33%	0.223%
25	21.231	3083	3088	3100	rBV4	661757	1146321	7.50%	1.259%
26	21.589	3140	3149	3154	rBV3	3172548	5996032	39.23%	6.588%
27	21.636	3154	3157	3165	rVB	684191	1093788	7.16%	1.202%
28	21.918	3199	3205	3213	rBV2	450513	875101	5.73%	0.961%
29	22.382	3279	3284	3291	rVB7	100369	201121	1.32%	0.221%
30	23.064	3396	3400	3407	rVB3	105513	207375	1.36%	0.228%
31	23.757	3509	3518	3524	rBV3	323195	1006946	6.59%	1.106%
32	23.857	3530	3535	3542	rVB2	427676	939455	6.15%	1.032%
33	24.462	3632	3638	3651	rVB2	233028	646686	4.23%	0.710%
34	24.609	3656	3663	3676	rVB2	159978	475831	3.11%	0.523%
35	24.762	3679	3689	3703	rBV	1768517	5532770	36.20%	6.079%
36	25.913	3874	3885	3896	rBV	1192498	3663811	23.97%	4.025%

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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_G\Methods\8270-BG122923.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

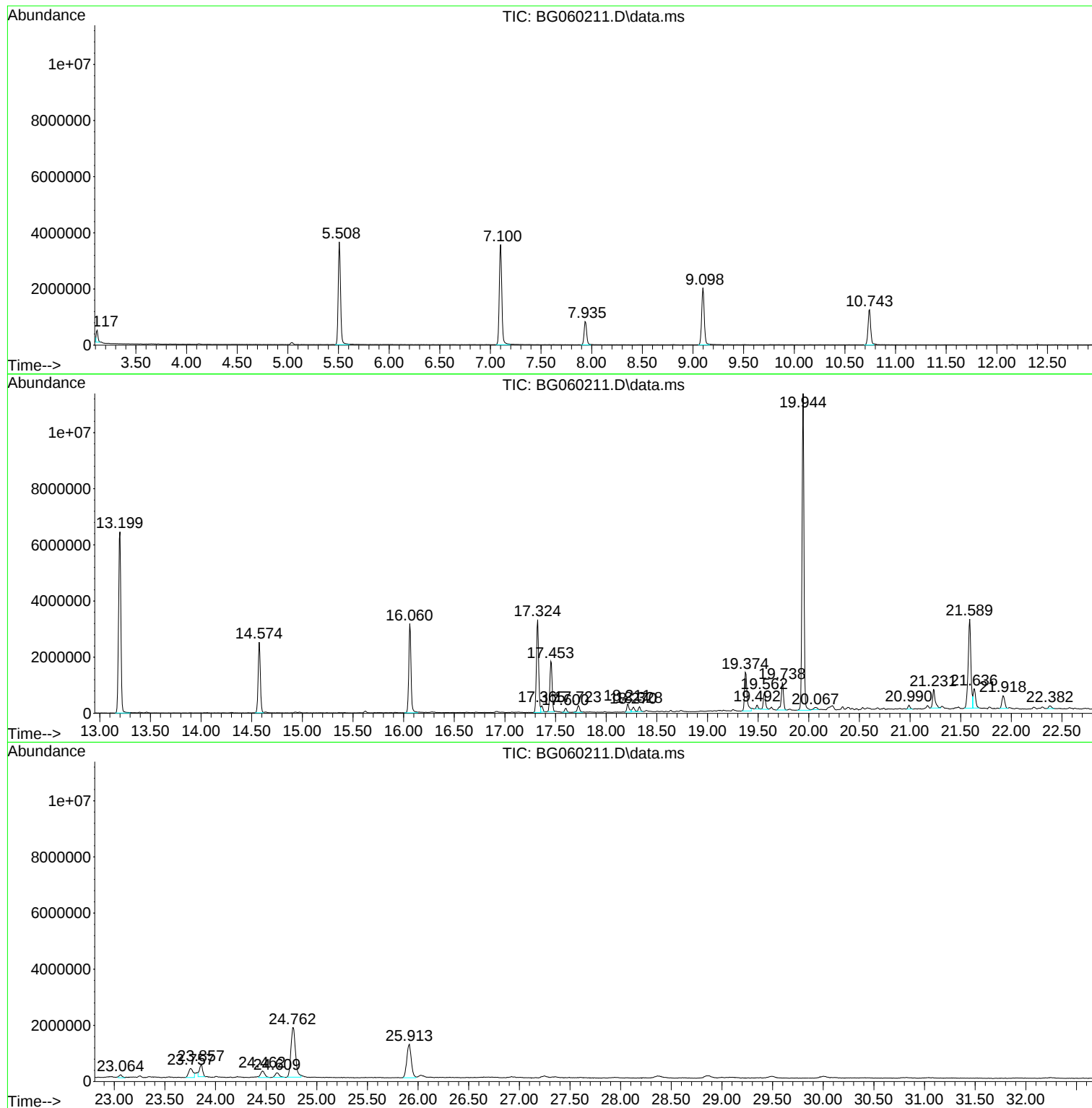
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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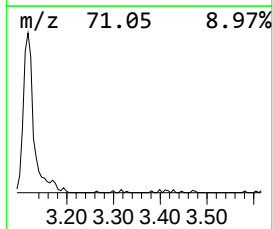
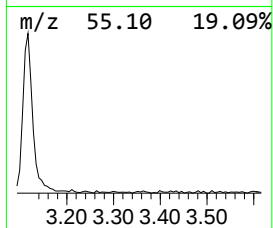
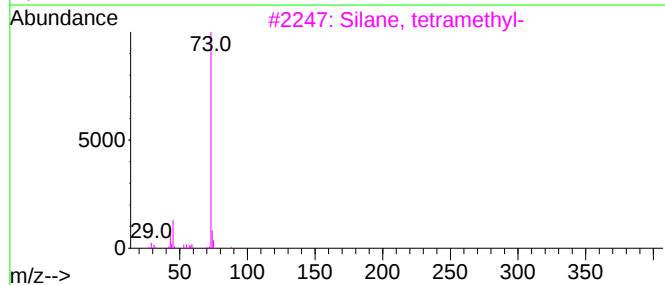
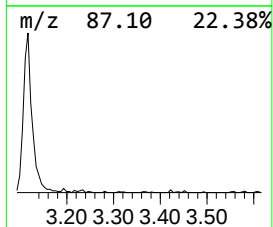
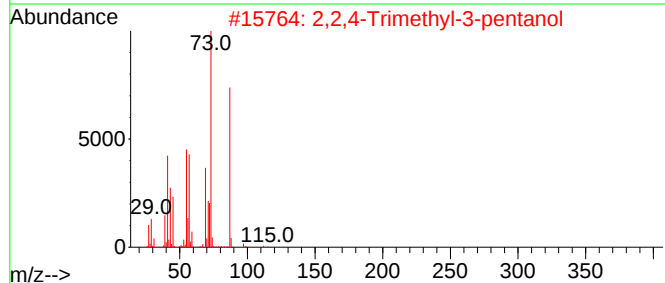
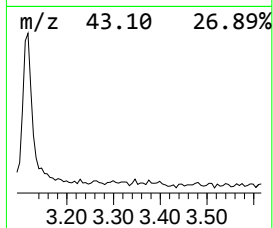
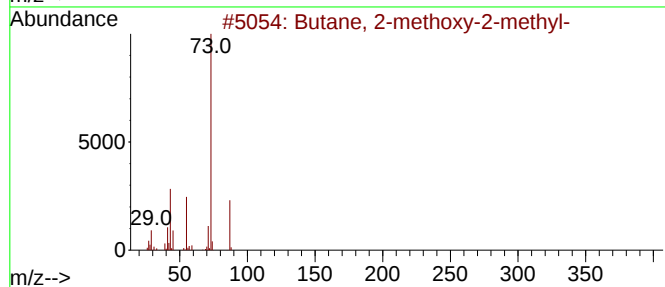
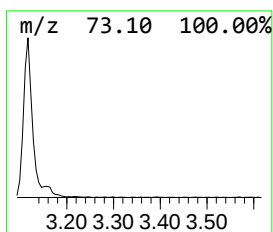
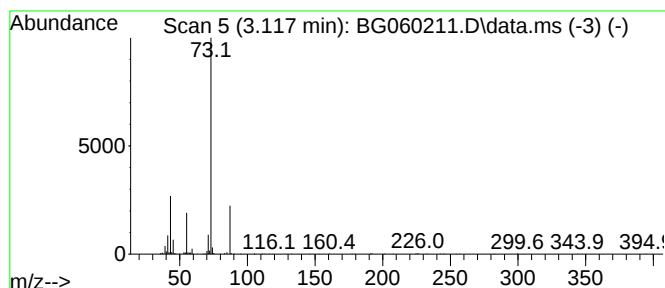
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.117	6.15 ng	443259	1,4-Dichlorobenzene-d4	7.935

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	50
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1-Propene-1-thiol	74	C3H6S	000925-89-3	9



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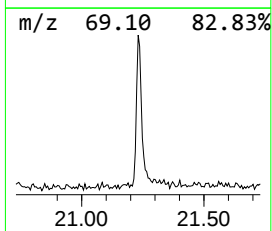
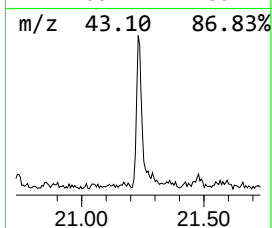
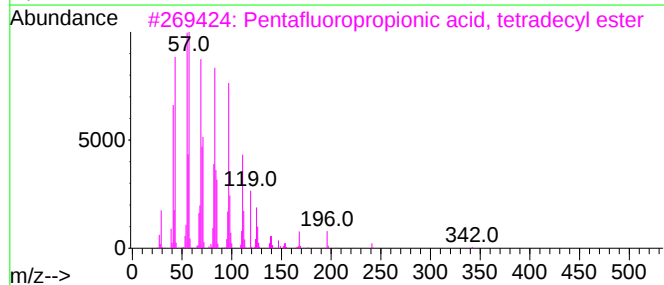
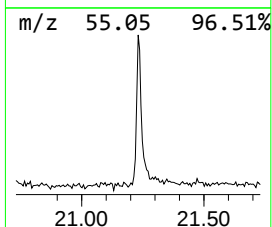
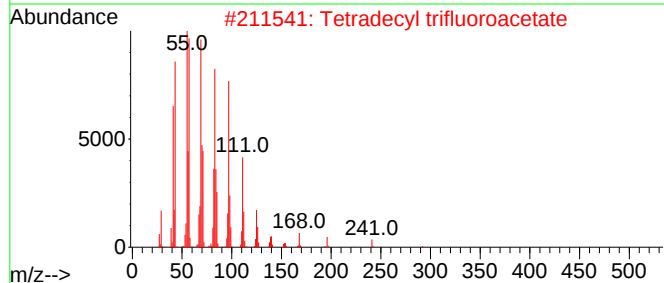
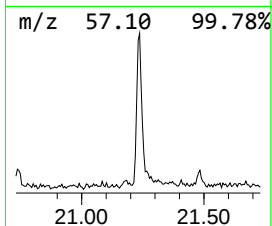
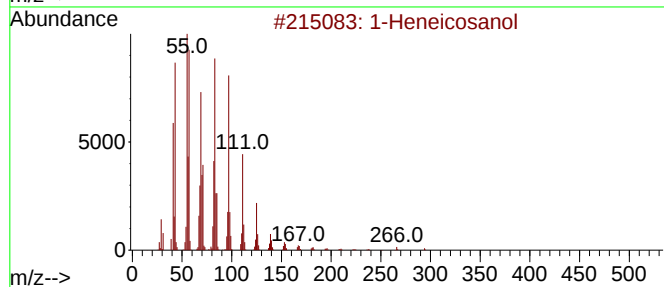
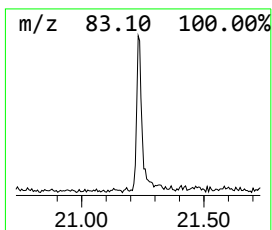
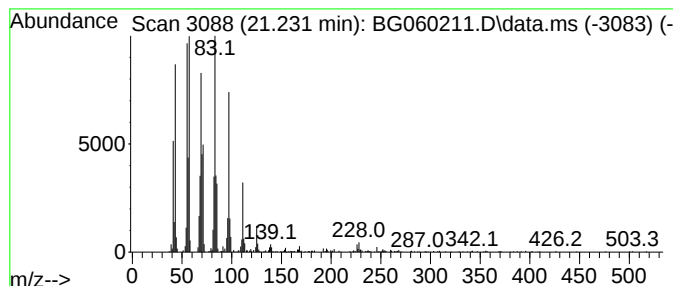
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.231	3.82 ng	1146320	Chrysene-d12	21.589

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	91
2			Tetradecyl trifluoroacetate	310	C16H29F3O2	006222-02-2	91
3			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	91
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5			Behenic alcohol	326	C22H46O	000661-19-8	91



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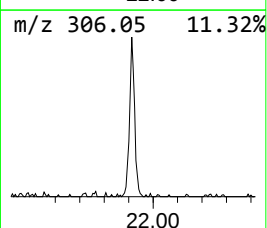
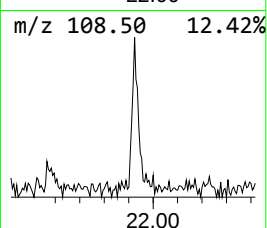
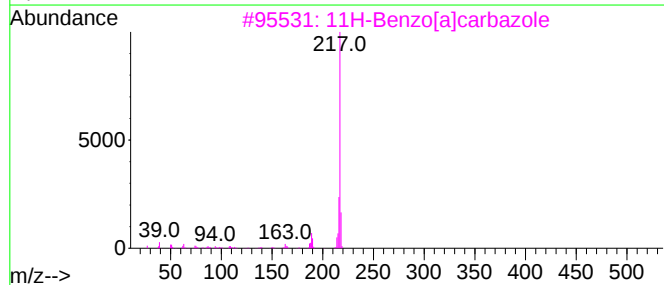
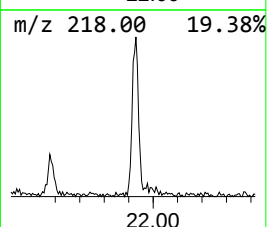
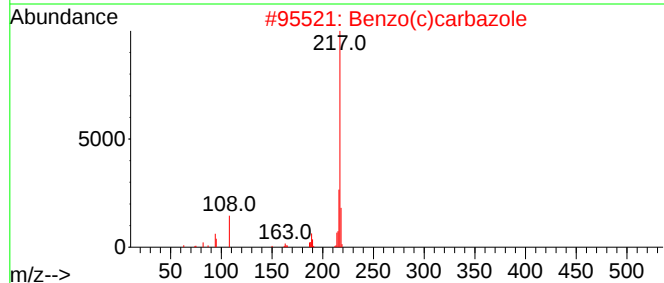
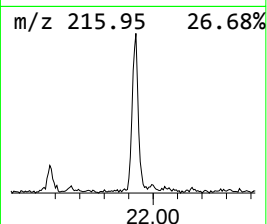
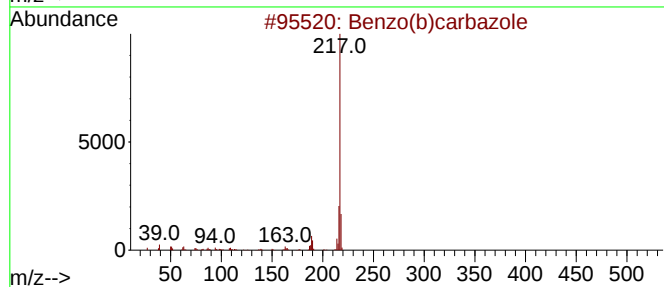
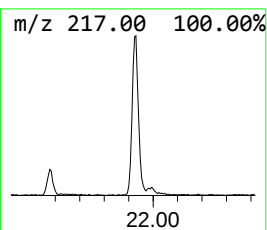
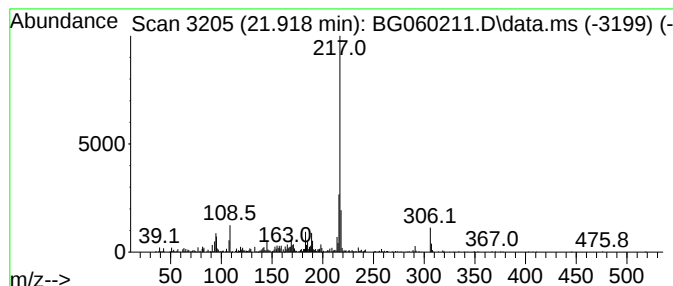
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzo(b)carbazole Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.918	2.92 ng	875101	Chrysene-d12	21.589

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo(b)carbazole	217	C16H11N	000243-28-7	96
2			Benzo(c)carbazole	217	C16H11N	034777-33-8	89
3			11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	86
4			7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	76
5			11H-Indeno(1,2-b)quinoline	217	C16H11N	000243-51-6	64



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ClientSampleId :
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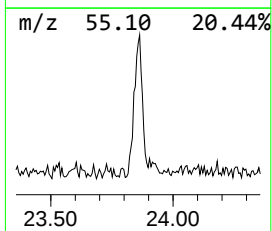
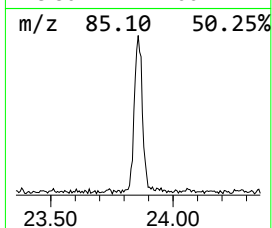
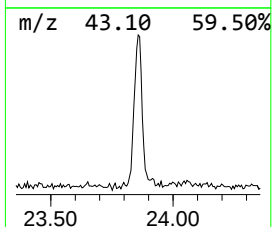
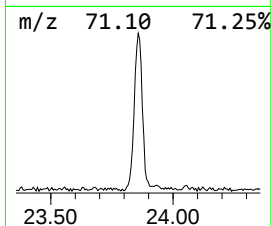
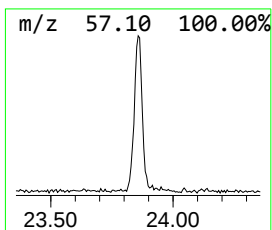
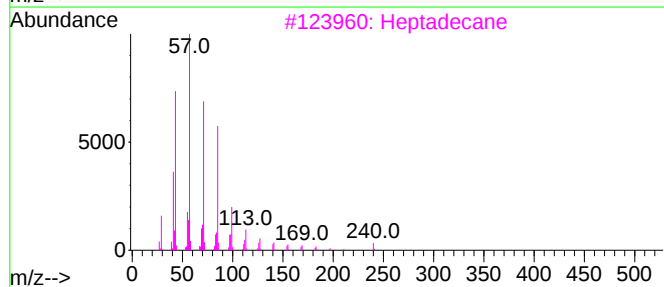
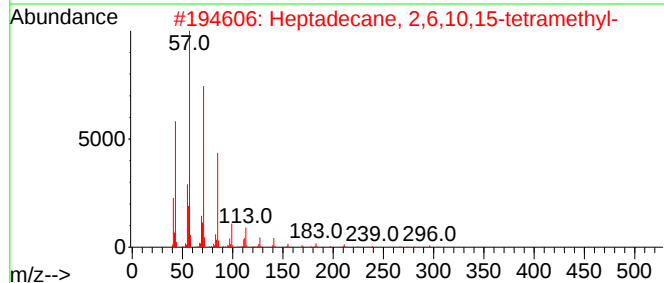
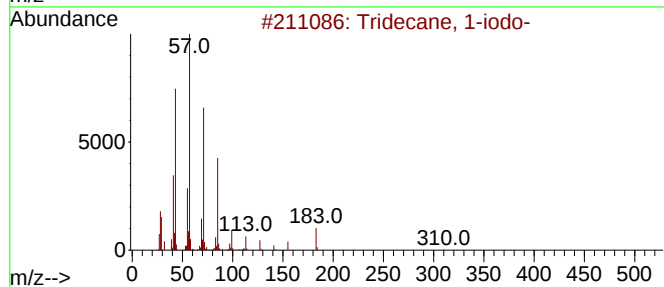
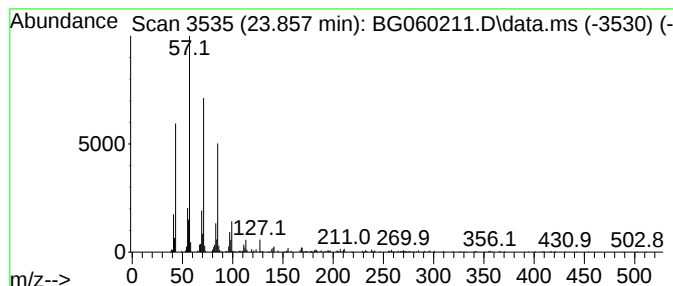
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Peak Number 4 Tridecane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.857	3.40 ng	939455	Perylene-d12	24.768

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
3			Heptadecane	240	C17H36	000629-78-7	86
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
5			Nonadecane	268	C19H40	000629-92-5	80



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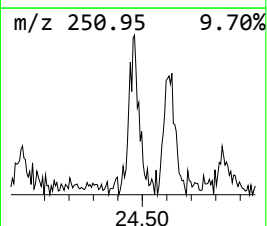
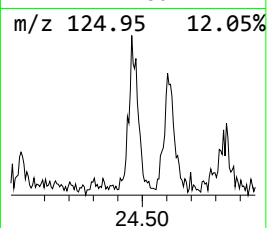
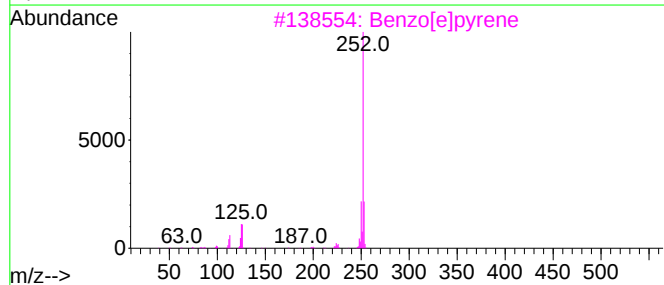
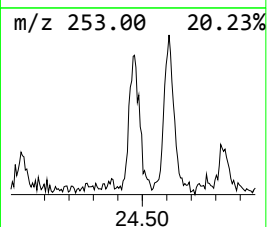
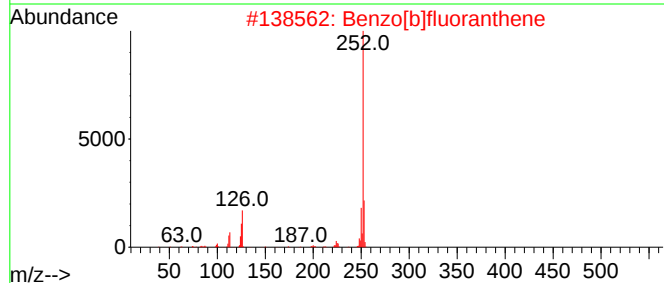
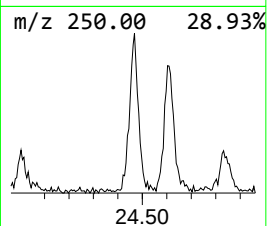
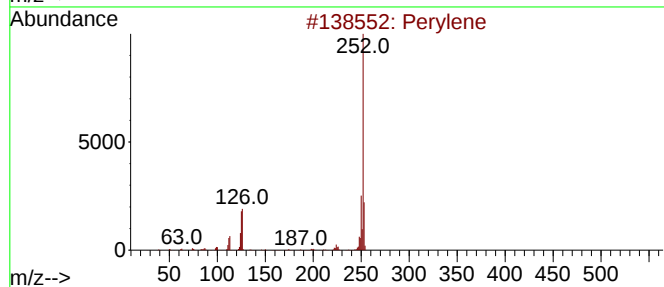
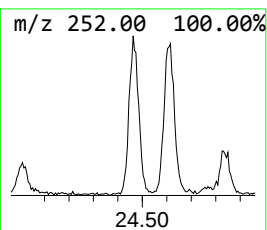
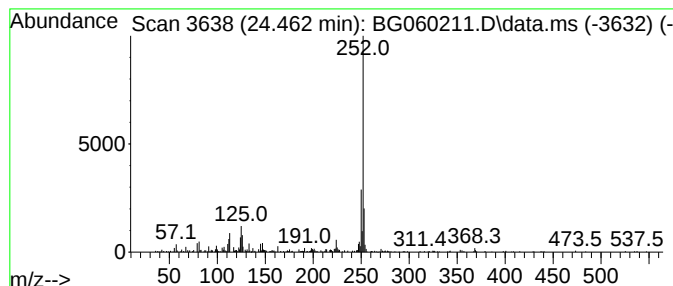
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Perylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.462	2.34 ng	646686	Perylene-d12	24.768

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Perylene	252	C20H12	000198-55-0	89
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	89
3			Benzo[e]pyrene	252	C20H12	000192-97-2	89
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	86
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010224\
Data File : BG060211.D
Acq On : 3 Jan 2024 3:16
Operator : MA/JU
Sample : 05899-24
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GHL-SS-19-2-4

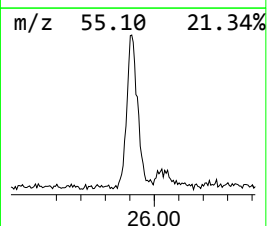
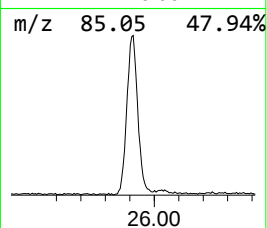
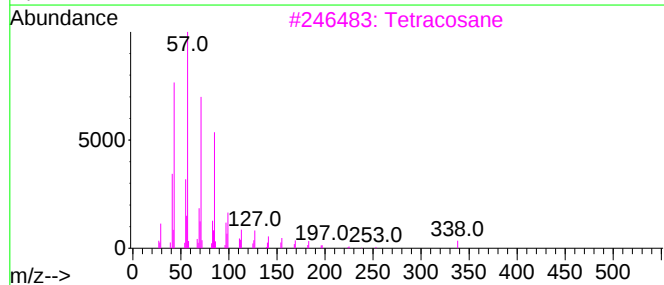
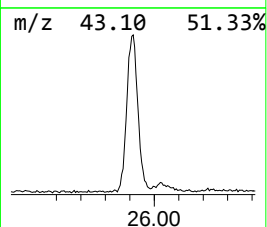
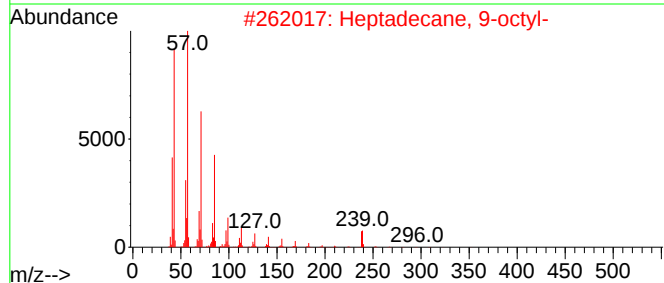
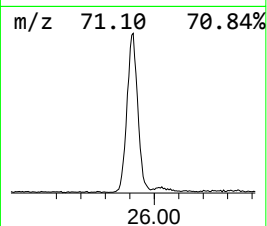
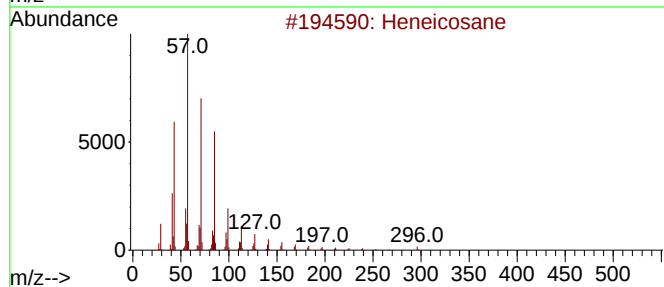
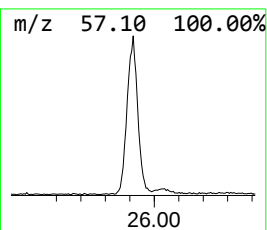
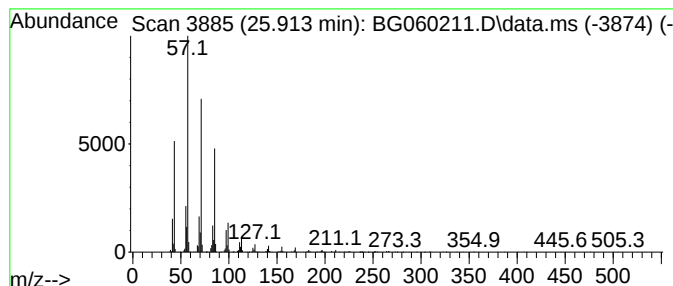
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heneicosane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.913	13.24 ng	3663810	Perylene-d12	24.768

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
3			Tetracosane	338	C24H50	000646-31-1	87
4			Tridecane, 7-propyl-	226	C16H34	055045-09-5	87
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010224\
Data File : BG060211.D
Acq On : 3 Jan 2024 3:16
Operator : MA/JU
Sample : 05899-24
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GHL-SS-19-2-4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.117	6.2	ng	443259	1	7.935	1441760	20.0
1-Heneicosanol	21.231	3.8	ng	1146320	5	21.589	5996030	20.0
Benzo(b)carbazole	21.918	2.9	ng	875101	5	21.589	5996030	20.0
Tridecane, 1-iodo-	23.857	3.4	ng	939455	6	24.768	5532770	20.0
Perylene	24.462	2.3	ng	646686	6	24.768	5532770	20.0
Heneicosane	25.913	13.2	ng	3663810	6	24.768	5532770	20.0