

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
 Data File : BP019813.D
 Acq On : 21 Feb 2024 18:02
 Operator : MA/JU
 Sample : P1537-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 5

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\8270E-BP013124.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP019813.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	21	rVB	692399	820108	31.20%	3.651%
2	4.069	161	167	175	rVB3	45260	75546	2.87%	0.336%
3	4.305	203	207	216	rBV2	34771	55136	2.10%	0.245%
4	4.452	228	232	240	rVB2	20411	29642	1.13%	0.132%
5	4.640	259	264	268	rBV	77910	113366	4.31%	0.505%
6	4.687	268	272	284	rVB3	37659	81804	3.11%	0.364%
7	5.016	321	328	337	rBV	254738	377464	14.36%	1.680%
8	5.469	398	405	414	rBV	966877	1475783	56.14%	6.570%
9	6.093	502	511	521	rBV2	23010	72355	2.75%	0.322%
10	6.716	607	617	624	rBV8	24898	75330	2.87%	0.335%
11	7.063	669	676	685	rBV	922187	1486342	56.54%	6.617%
12	7.593	757	766	775	rBV10	14611	32771	1.25%	0.146%
13	7.893	810	817	826	rVB	326473	527289	20.06%	2.347%
14	8.863	976	982	987	rBV6	15344	34434	1.31%	0.153%
15	9.063	1008	1016	1026	rBV	587860	1010109	38.43%	4.497%
16	9.357	1058	1066	1070	rBV9	9345	26486	1.01%	0.118%
17	10.051	1177	1184	1195	rBV6	20789	66607	2.53%	0.297%
18	10.693	1282	1293	1305	rBV	384227	690467	26.27%	3.074%
19	13.151	1704	1711	1726	rBV	1395993	2215474	84.28%	9.862%
20	14.251	1893	1898	1907	rVB	25351	41361	1.57%	0.184%
21	14.422	1919	1927	1933	rBV7	8108	26659	1.01%	0.119%
22	14.528	1938	1945	1952	rVV	523103	826859	31.46%	3.681%
23	14.587	1952	1955	1965	rVB	19827	38792	1.48%	0.173%
24	15.581	2119	2124	2129	rBV4	15011	26382	1.00%	0.117%
25	16.022	2190	2199	2221	rBV	1052005	1742477	66.29%	7.757%
26	16.257	2233	2239	2248	rVB	12901	29887	1.14%	0.133%
27	16.586	2288	2295	2300	rBV4	13606	30869	1.17%	0.137%
28	17.286	2406	2414	2418	rBV	585944	911372	34.67%	4.057%
29	17.328	2418	2421	2429	rVV	224508	327867	12.47%	1.460%
30	17.416	2431	2436	2441	rVB	63457	99442	3.78%	0.443%
31	17.869	2509	2513	2521	rVB5	14605	26804	1.02%	0.119%
32	18.180	2562	2566	2575	rVB2	215445	407352	15.50%	1.813%
33	18.280	2579	2583	2587	rVV	32482	44394	1.69%	0.198%
34	18.339	2588	2593	2597	rVV2	88630	165589	6.30%	0.737%
35	18.375	2597	2599	2605	rVB2	50391	62494	2.38%	0.278%
36	18.639	2640	2644	2648	rBV	33720	46554	1.77%	0.207%

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ClientSampleId :
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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

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37	18.686	2648	2652	2660	rVB5	20541	38425	1.46%	0.171%
38	18.922	2689	2692	2695	rBV	21743	26333	1.00%	0.117%
39	19.039	2708	2712	2716	rBV	58052	85611	3.26%	0.381%
40	19.104	2716	2723	2725	rVV5	26568	49899	1.90%	0.222%
41	19.180	2731	2736	2743	rBV5	32238	65041	2.47%	0.290%
42	19.298	2750	2756	2762	rBV	473688	665698	25.32%	2.963%
43	19.416	2769	2776	2785	rBV5	98554	177659	6.76%	0.791%
44	19.622	2806	2811	2812	rBV2	90957	124991	4.75%	0.556%
45	19.645	2812	2815	2824	rVV	435194	623901	23.73%	2.777%
46	19.722	2824	2828	2833	rVB3	41572	59838	2.28%	0.266%
47	19.851	2844	2850	2862	rVB	2171228	2628628	100.00%	11.702%
48	19.969	2865	2870	2876	rBV3	42598	103059	3.92%	0.459%
49	20.098	2888	2892	2894	rBV4	38561	64651	2.46%	0.288%
50	20.133	2895	2898	2905	rVB	100386	132098	5.03%	0.588%
51	20.227	2911	2914	2918	rBV	58772	76444	2.91%	0.340%
52	20.286	2920	2924	2928	rVB3	61967	92878	3.53%	0.413%
53	20.421	2944	2947	2951	rBV	43736	51703	1.97%	0.230%
54	20.469	2952	2955	2958	rVB	42155	52944	2.01%	0.236%
55	21.098	3059	3062	3068	rBV3	170346	219236	8.34%	0.976%
56	21.374	3103	3109	3113	rBV2	802542	1403558	53.40%	6.248%
57	21.416	3113	3116	3124	rVB	235936	332661	12.66%	1.481%
58	23.057	3391	3395	3400	rBV	127404	282300	10.74%	1.257%
59	23.792	3514	3520	3527	rBV	497021	984677	37.46%	4.383%

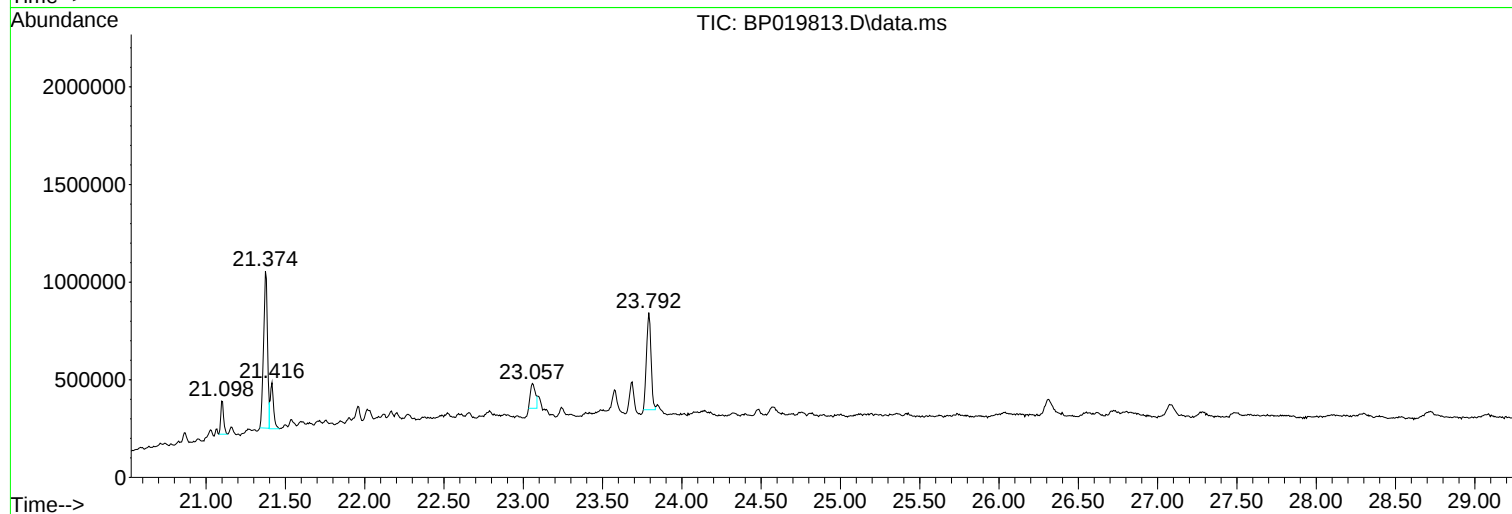
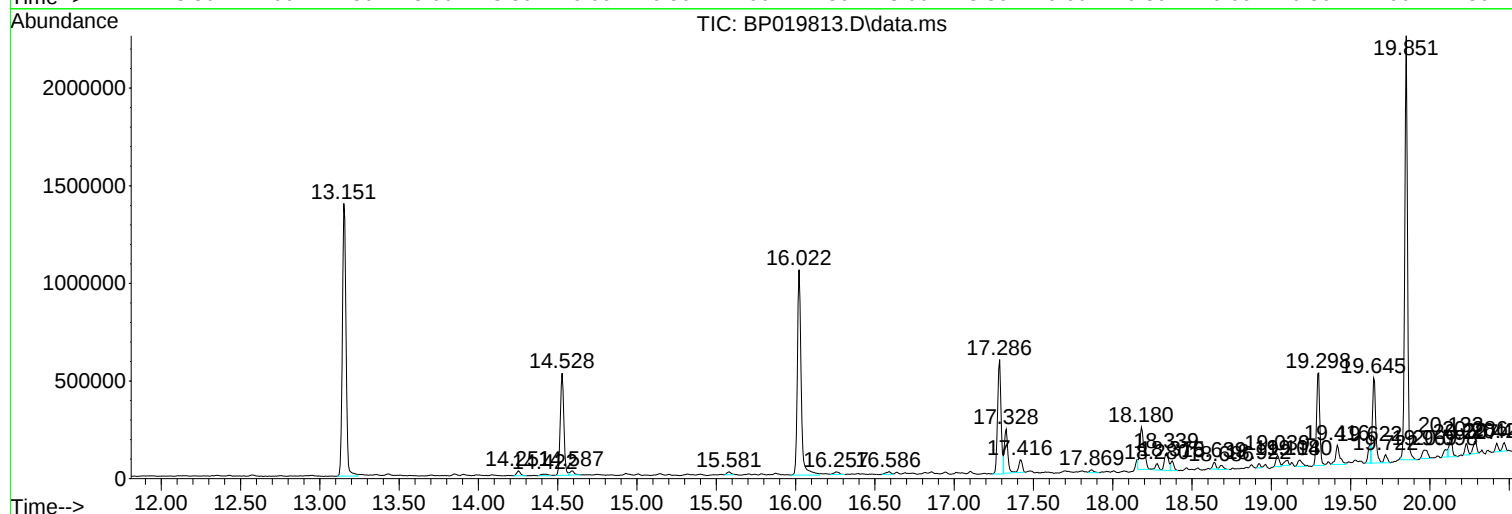
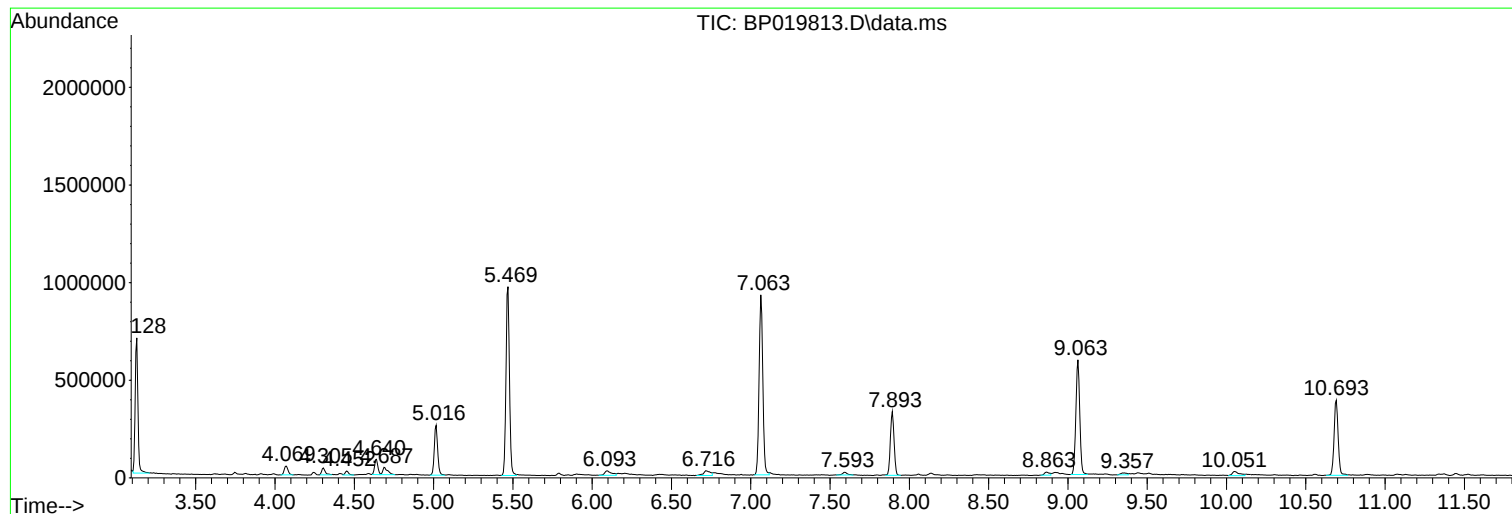
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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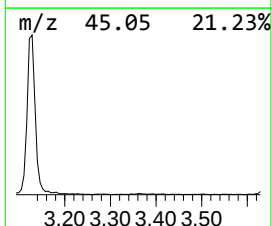
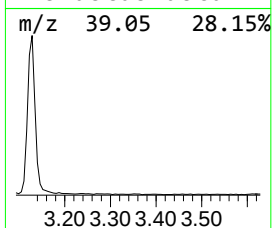
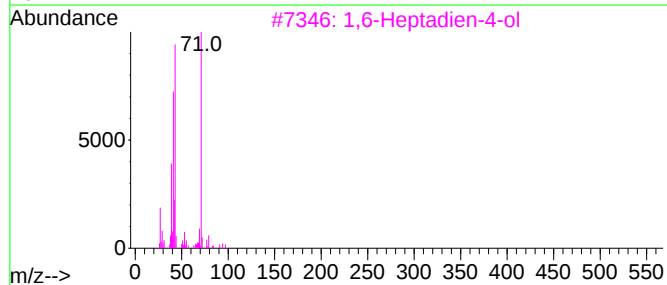
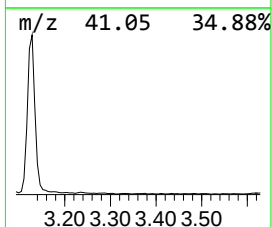
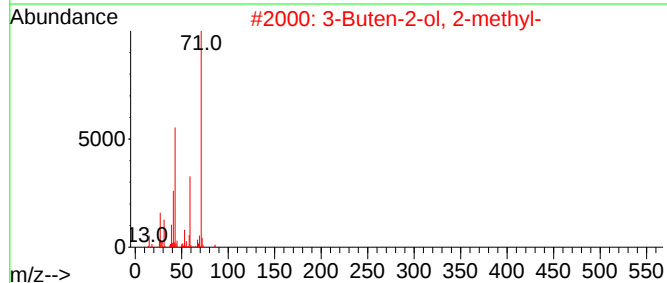
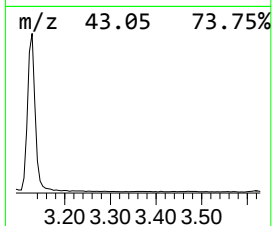
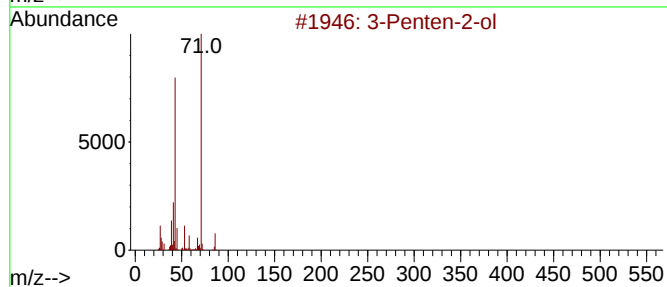
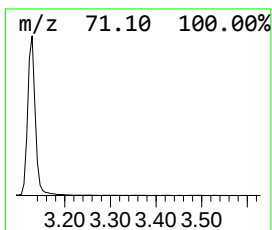
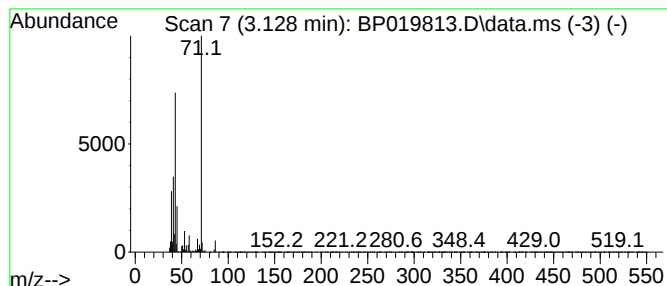
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-ol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	31.11 ng	820108	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-ol	86	C5H10O	001569-50-2	83
2			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	64
3			1,6-Heptadien-4-ol	112	C7H12O	002883-45-6	59
4			2-Furanmethanol, tetrahydro-, ac...	144	C7H12O3	000637-64-9	53
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	50



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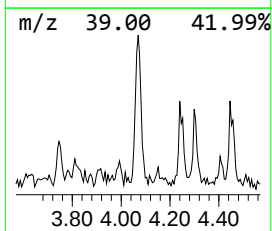
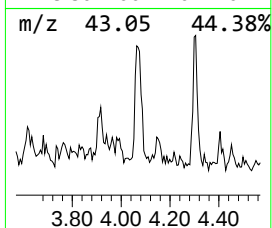
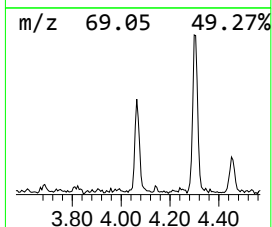
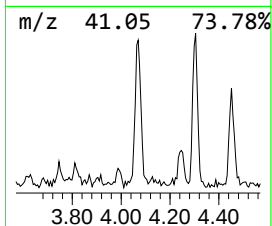
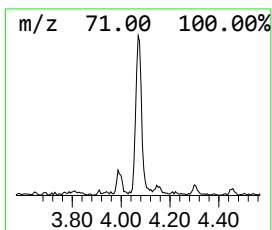
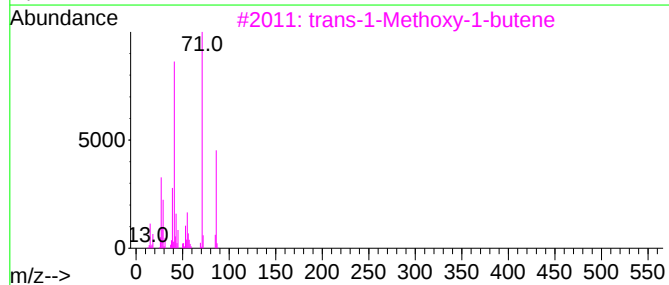
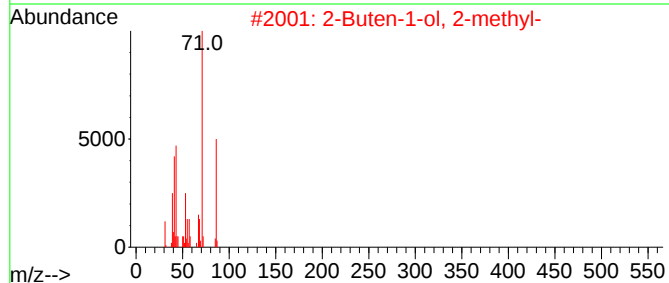
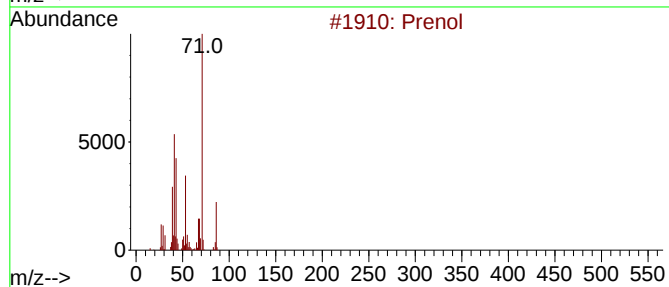
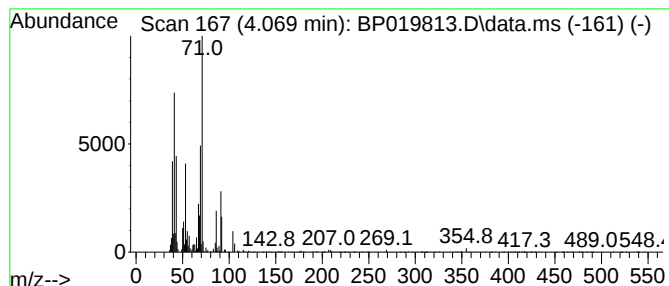
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Prenol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.069	2.87 ng	75546	1,4-Dichlorobenzene-d4	7.893

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Prenol	86	C5H10O	000556-82-1	81
2		2-Buten-1-ol, 2-methyl-	86	C5H10O	004675-87-0	64
3		trans-1-Methoxy-1-butene	86	C5H10O	010034-13-6	18
4		2-Butene, 1-methoxy-, (E)-	86	C5H10O	010034-14-7	16
5		1-Propene, 1-methoxy-2-methyl-	86	C5H10O	017574-84-4	14



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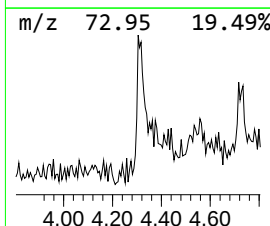
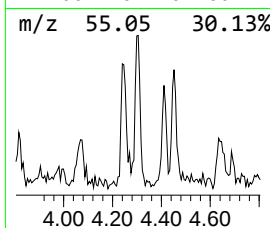
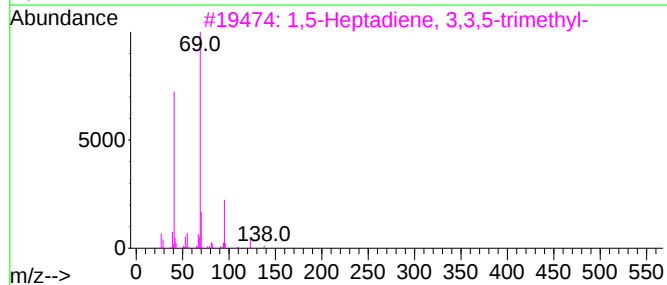
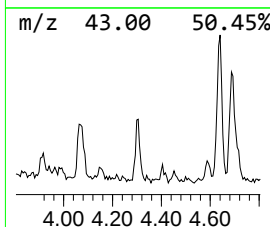
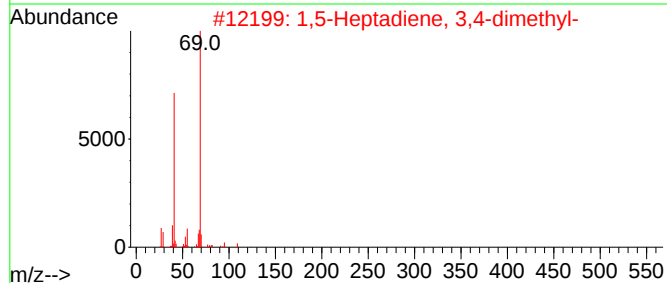
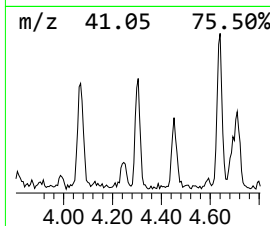
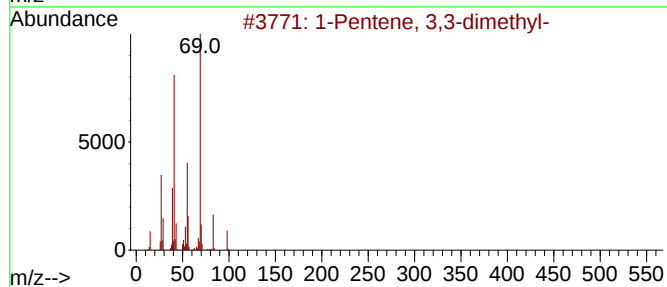
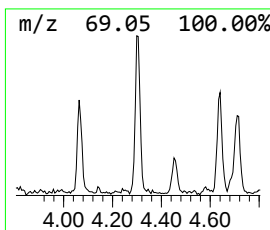
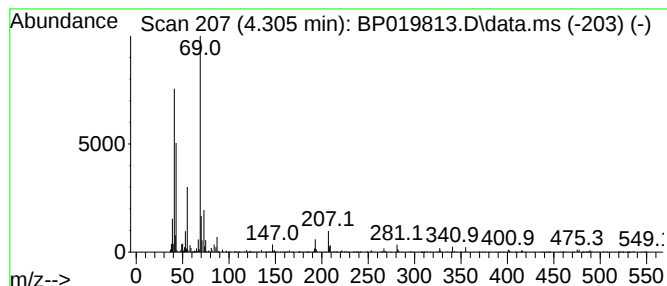
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 1-Pentene, 3,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.305	2.09 ng	55136	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	50
2			1,5-Heptadiene, 3,4-dimethyl-	124	C9H16	1000061-77-9	50
3			1,5-Heptadiene, 3,3,5-trimethyl-	138	C10H18	074630-29-8	33
4			4-Trifluoroacetoxyoctane	226	C10H17F3O2	116465-17-9	33
5			3-Butyn-2-ol, 2-methyl-	84	C5H8O	000115-19-5	33



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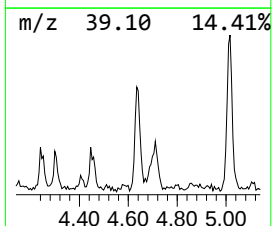
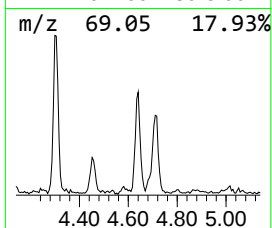
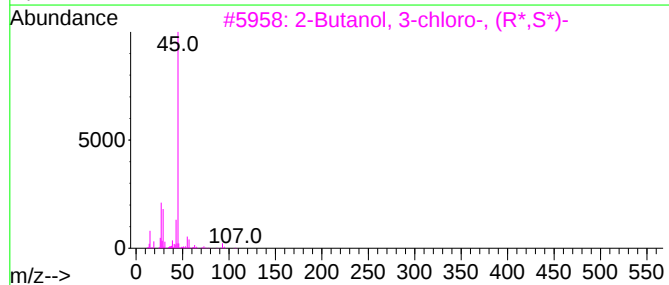
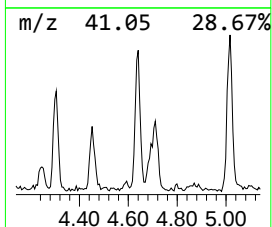
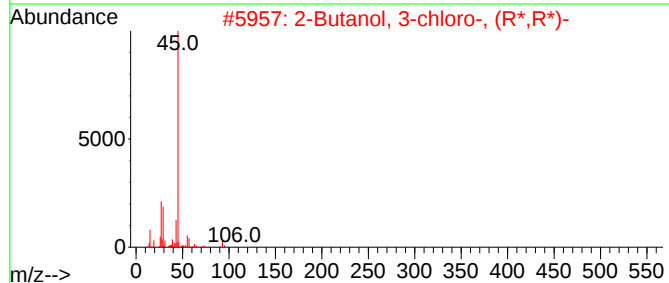
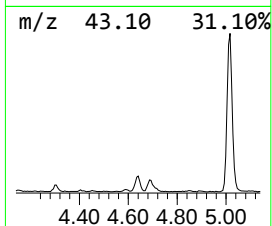
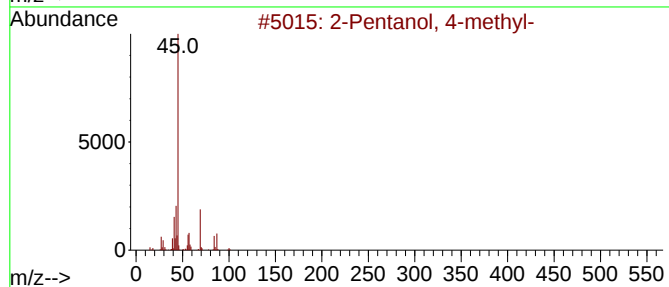
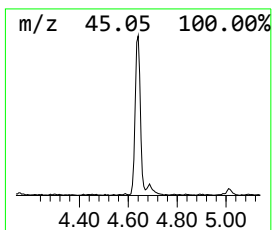
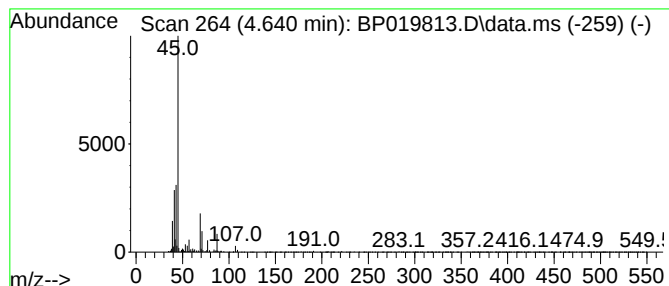
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2-Pentanol, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.640	4.30 ng	113366	1,4-Dichlorobenzene-d4	7.893

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	64
2		2-Butanol, 3-chloro-, (R*,R*)-	108	C4H9ClO	010325-40-3	42
3		2-Butanol, 3-chloro-, (R*,S*)-	108	C4H9ClO	010325-41-4	42
4		Formic acid, 1-methylethyl ester	88	C4H8O2	000625-55-8	40
5		2-Hexanol	102	C6H14O	000626-93-7	39



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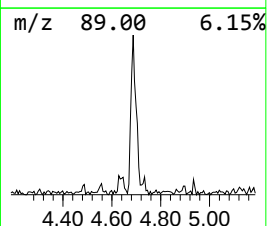
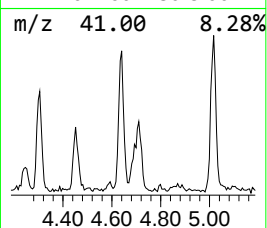
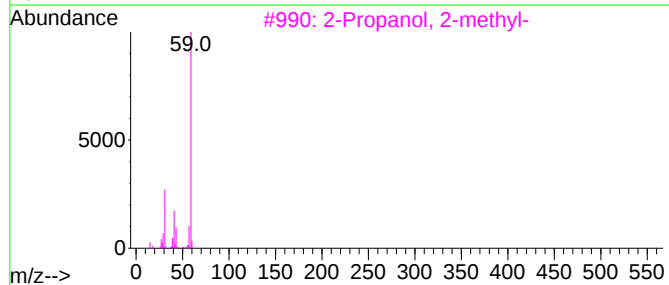
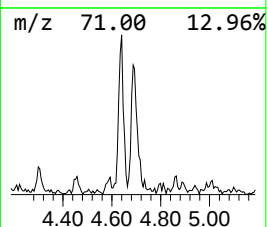
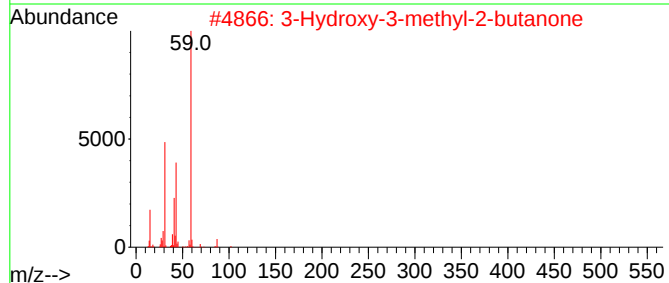
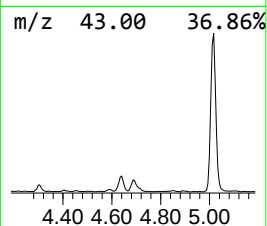
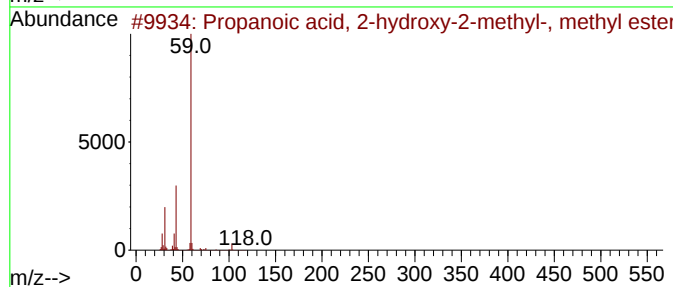
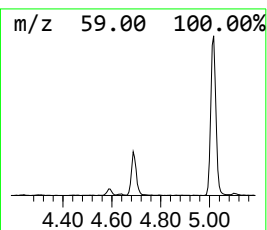
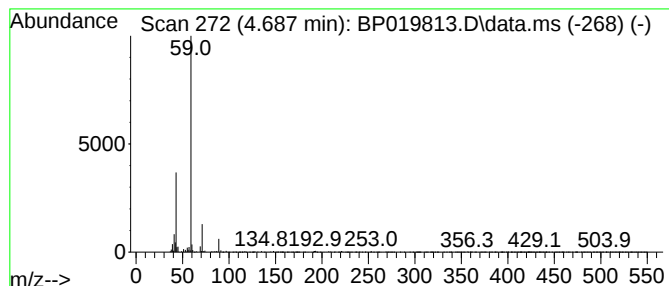
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown4.687 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.687	3.10 ng	81804	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-hydroxy-2-meth...	118	C5H10O3	002110-78-3	45
2			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	38
3			2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	38
4			Guanidine	59	CH5N3	000113-00-8	36
5			Butanamide, 2-hydroxy-2,N-dimethyl-	117	C5H11NO2	039961-72-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
Operator : MA/JU
Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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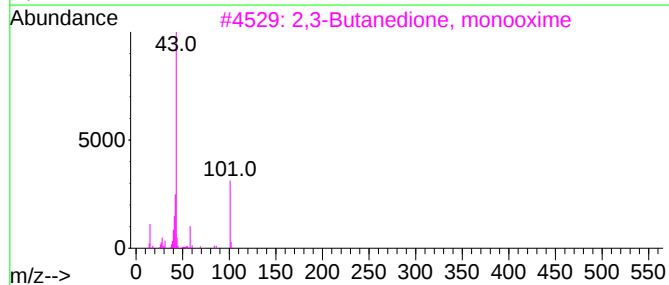
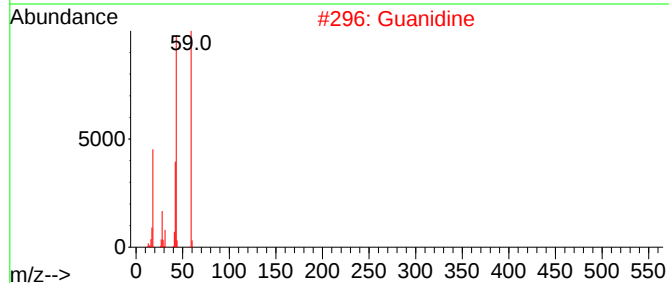
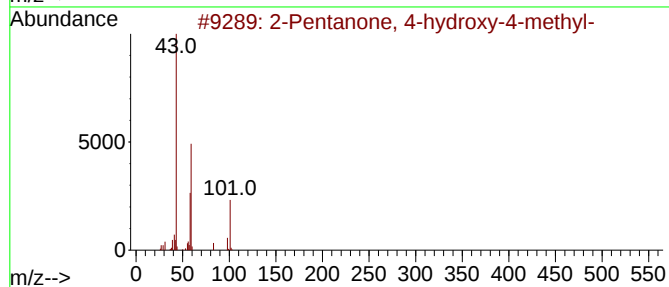
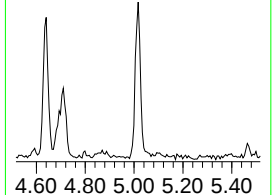
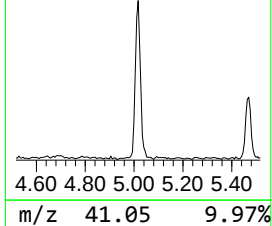
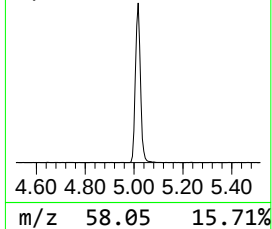
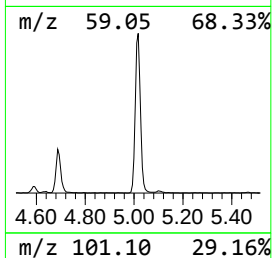
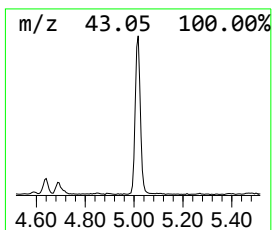
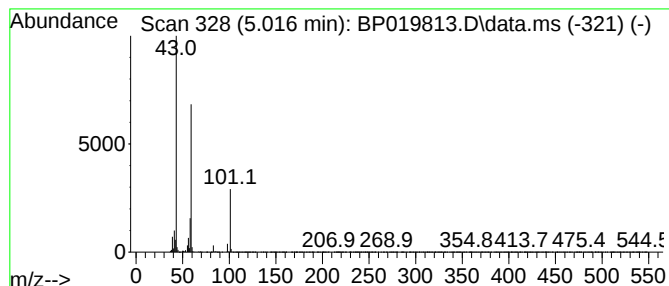
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	14.32 ng	377464	1,4-Dichlorobenzene-d4	7.893

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Guanidine	59	CH5N3	000113-00-8	43
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
Operator : MA/JU
Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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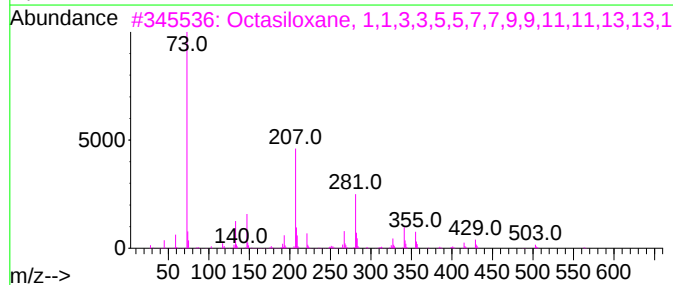
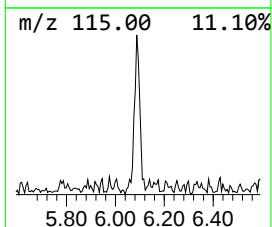
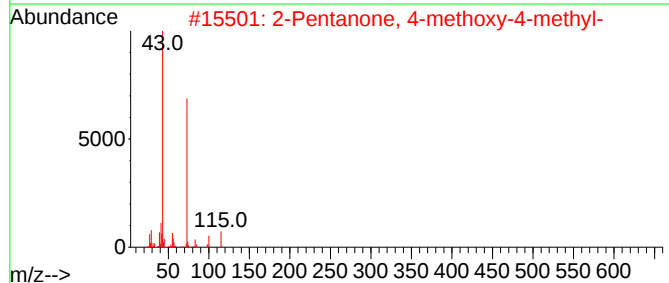
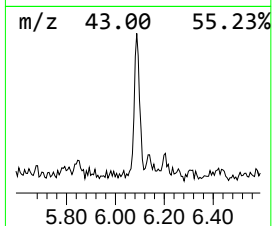
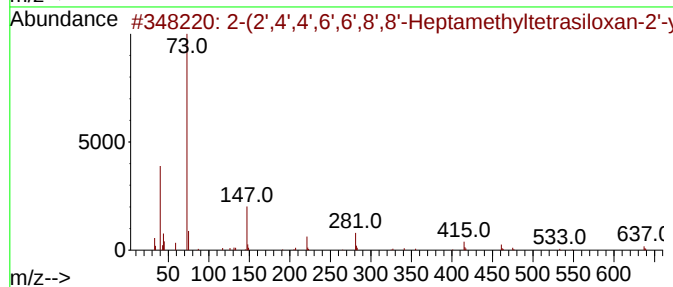
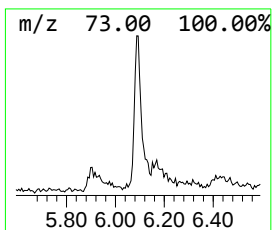
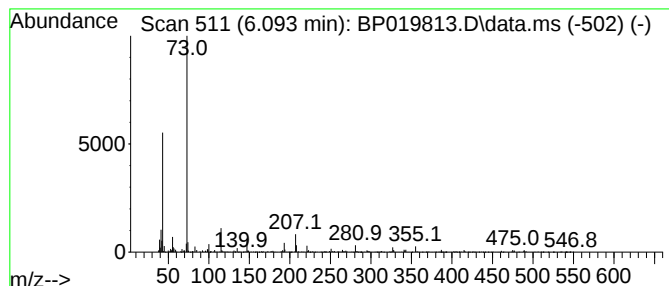
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.093 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.093	2.74 ng	72355	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2',4',4',6',6',8',8'-Heptamet...	652	C16H48010Si9	145344-72-5	37		
2	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	35		
3	Octasiloxane, 1,1,3,3,5,5,7,7,9,...	578	C16H5007Si8	019095-24-0	35		
4	Silane, (2-ethyl-5,5-dimethyl-4-...	208	C13H24Si	095798-05-3	25		
5	(S)-(-)-1,1'-Bi-2-naphthol, 0,0'...	430	C26H3002Si2	1000395-55-8	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
Operator : MA/JU
Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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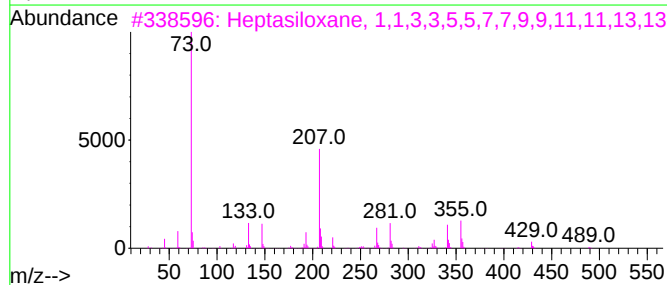
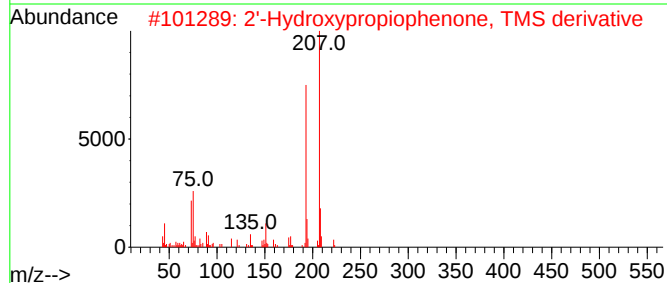
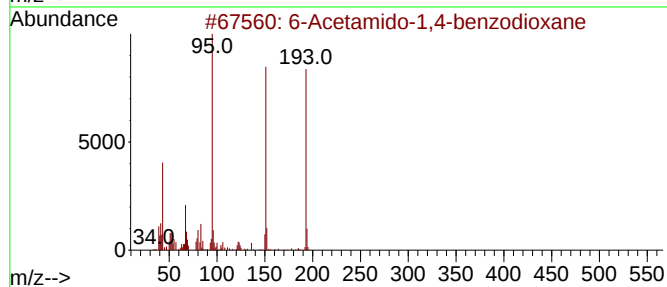
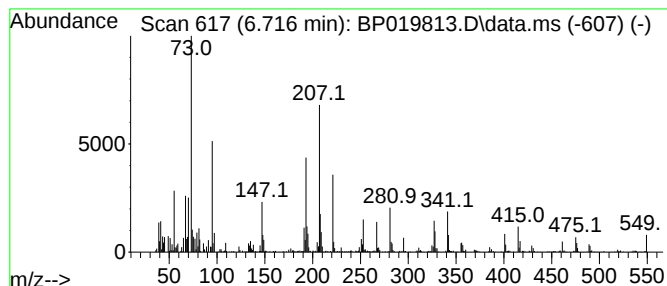
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown6.716 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.716	2.86 ng	75330	1,4-Dichlorobenzene-d4	7.893

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	35		
2	2'-Hydroxypropiophenone, TMS der...	222	C12H18O2Si	033342-87-9	25		
3	Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H44O6Si7	019095-23-9	22		
4	1-Pentamethyldisilanyloxy-4-pent...	354	C16H34OSi4	1000427-15-7	22		
5	Ethane, 1-(4,4,4-trifluoro-1,3-d...	308	C5H6F6S4	1000226-87-3	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
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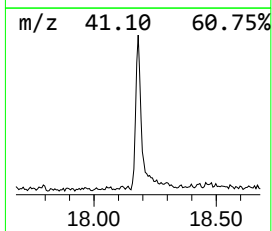
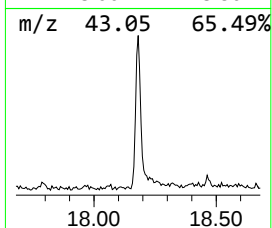
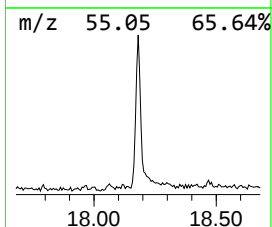
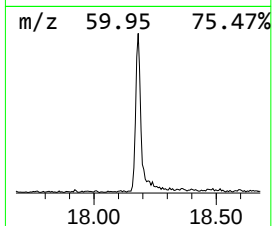
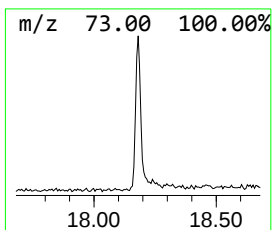
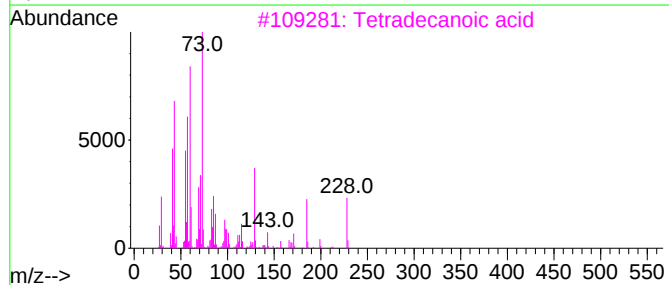
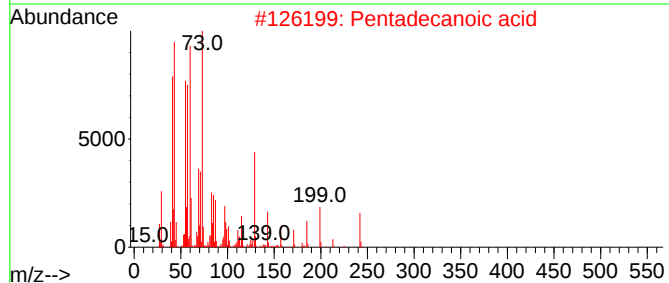
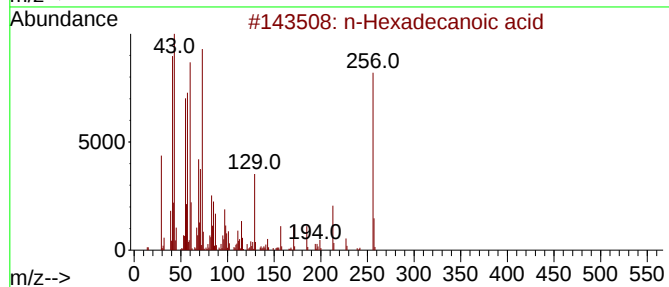
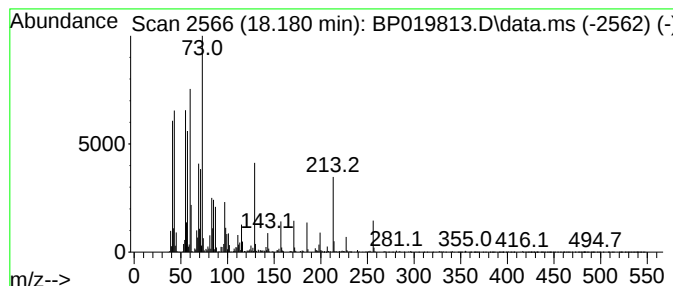
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.180	8.94 ng	407352	Phenanthrene-d10	17.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	81
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
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Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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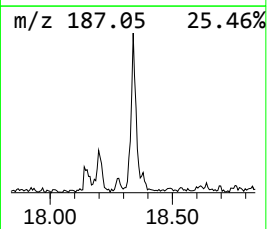
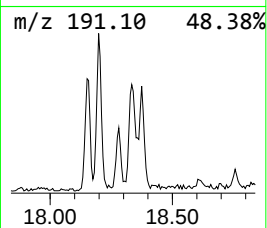
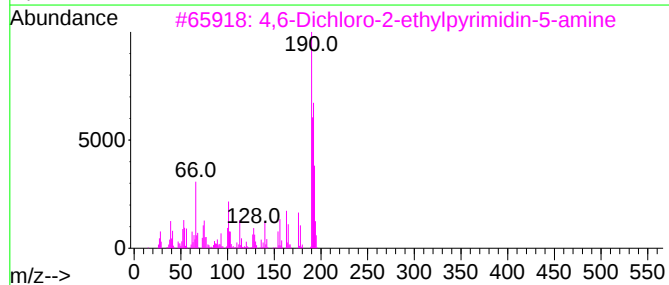
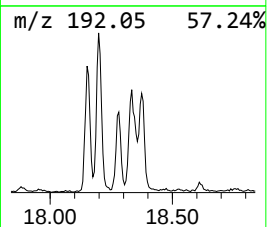
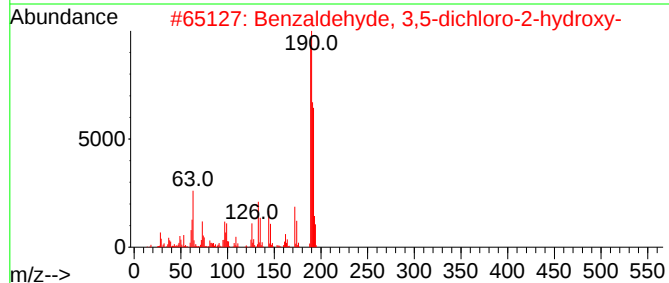
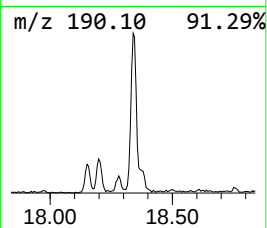
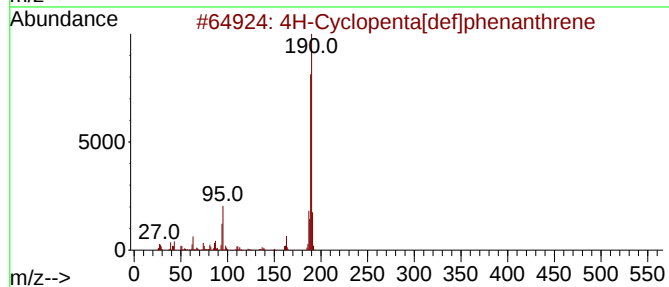
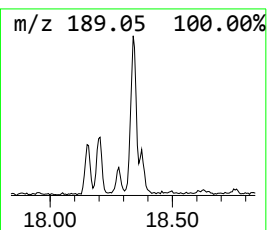
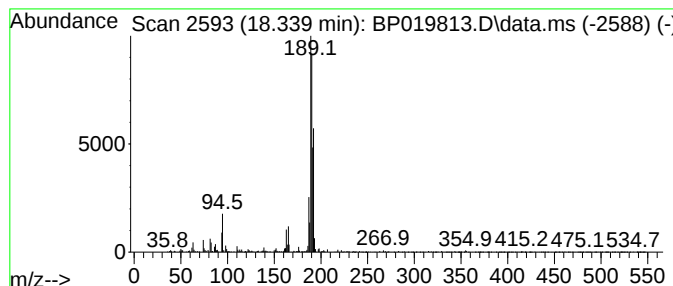
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.339	3.63 ng	165589	Phenanthrene-d10	17.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	47
3		4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	43
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
Operator : MA/JU
Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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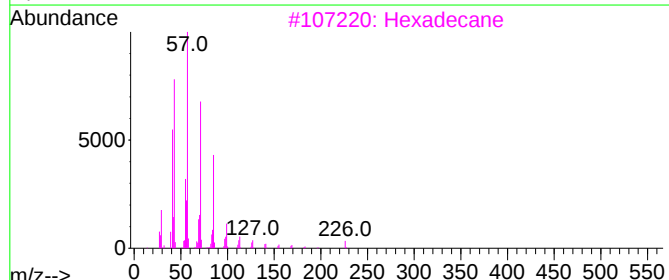
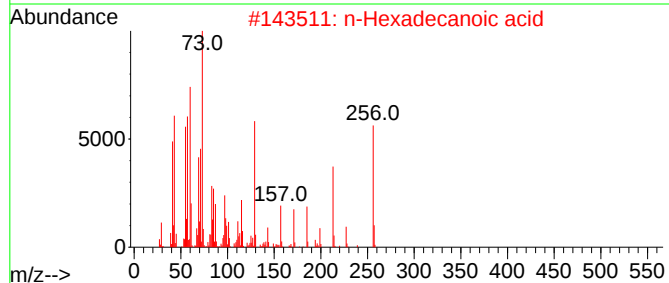
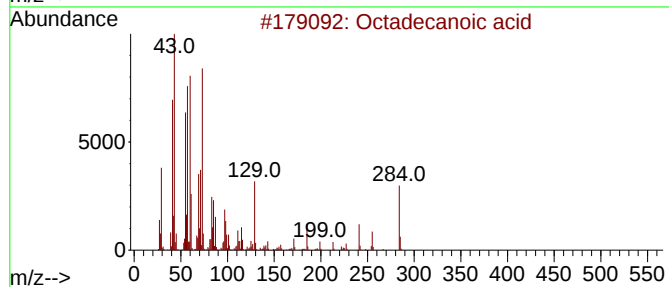
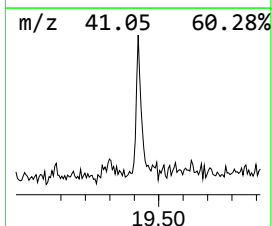
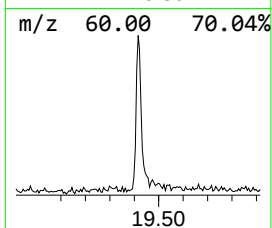
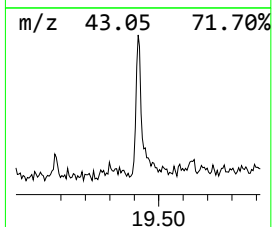
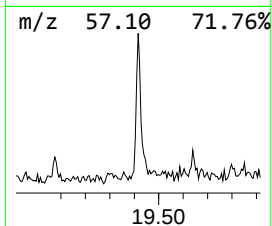
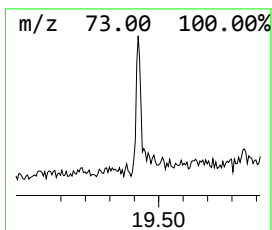
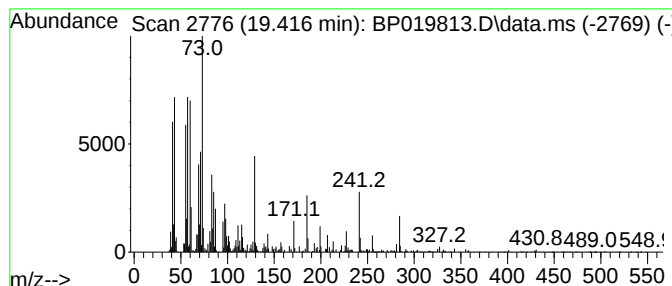
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.416	2.53 ng	177659	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3			Hexadecane	226	C16H34	000544-76-3	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
5			n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
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Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

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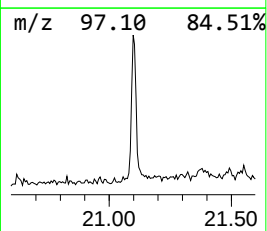
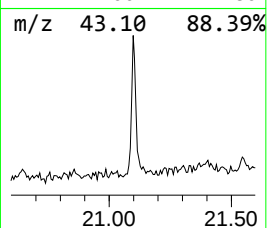
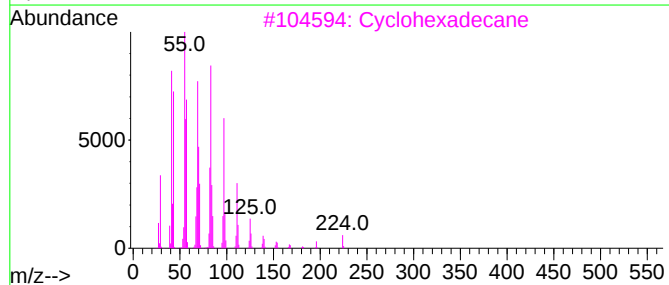
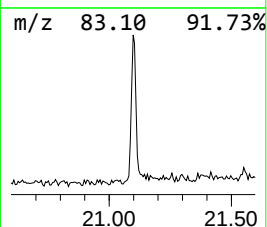
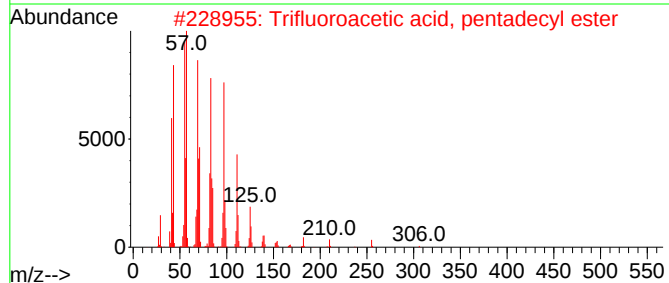
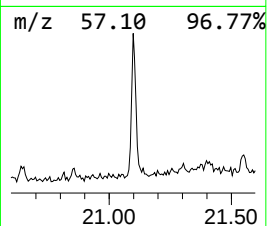
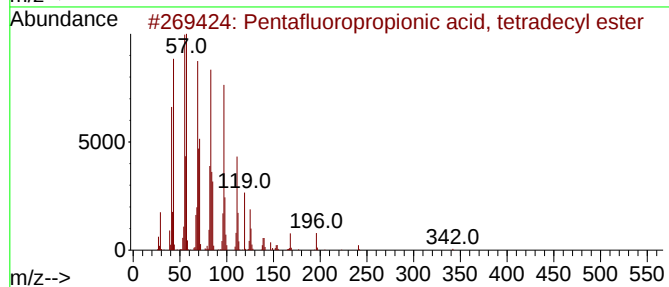
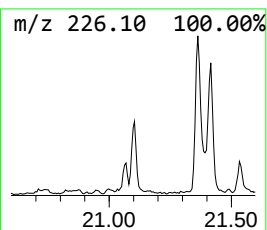
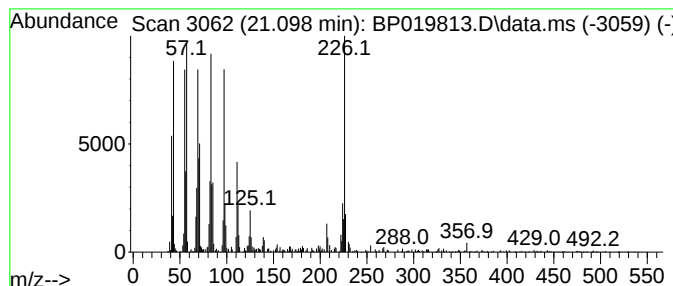
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentafluoropropionic acid, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	3.12 ng	219236	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	89
2			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	89
3			Cyclohexadecane	224	C16H32	000295-65-8	87
4			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	87
5			1-Hexacosene	364	C26H52	018835-33-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022124\
Data File : BP019813.D
Acq On : 21 Feb 2024 18:02
Operator : MA/JU
Sample : P1537-05
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-ol	3.128	31.1	ng	820108	1	7.893	527289	20.0
Prenol	4.069	2.9	ng	75546	1	7.893	527289	20.0
1-Pentene, 3,3-...	4.305	2.1	ng	55136	1	7.893	527289	20.0
2-Pentanol, 4-m...	4.640	4.3	ng	113366	1	7.893	527289	20.0
unknown4.687	4.687	3.1	ng	81804	1	7.893	527289	20.0
2-Pentanone, 4-...	5.016	14.3	ng	377464	1	7.893	527289	20.0
unknown6.093	6.093	2.7	ng	72355	1	7.893	527289	20.0
unknown6.716	6.716	2.9	ng	75330	1	7.893	527289	20.0
n-Hexadecanoic ...	18.180	8.9	ng	407352	4	17.286	911372	20.0
4H-Cyclopenta[d]...	18.339	3.6	ng	165589	4	17.286	911372	20.0
Octadecanoic acid	19.416	2.5	ng	177659	5	21.380	1403560	20.0
Pentafluoroprop...	21.098	3.1	ng	219236	5	21.380	1403560	20.0