

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
 Data File : BP005127.D
 Acq On : 31 Mar 2021 15:37
 Operator : CG/JU
 Sample : M1836-16
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B7-10-12

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-BP032321.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005127.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.699	97	104	111	rBV3	8927	14014	1.17%	0.195%
2	3.793	114	120	125	rBV	11126	17019	1.43%	0.237%
3	3.940	138	145	152	rBV3	8188	12956	1.08%	0.180%
4	5.311	372	378	384	rVV	429327	599720	50.22%	8.352%
5	5.369	384	388	398	rVB	22490	48240	4.04%	0.672%
6	5.740	441	451	457	rVB	15699	25054	2.10%	0.349%
7	6.699	607	614	621	rBV3	14188	25692	2.15%	0.358%
8	6.805	627	632	637	rVB	19333	27571	2.31%	0.384%
9	6.893	637	647	661	rBV	406507	679481	56.89%	9.463%
10	7.099	677	682	689	rVB	8429	13737	1.15%	0.191%
11	7.358	717	726	732	rVB	48810	80321	6.73%	1.119%
12	7.711	777	786	795	rBV	107163	176287	14.76%	2.455%
13	7.828	801	806	814	rVB2	13996	24205	2.03%	0.337%
14	8.087	844	850	855	rVB2	10929	17932	1.50%	0.250%
15	8.252	873	878	884	rBV3	9936	17265	1.45%	0.240%
16	8.346	889	894	909	rVB2	16509	32872	2.75%	0.458%
17	8.810	968	973	977	rBV	15460	23892	2.00%	0.333%
18	8.869	977	983	992	rVB	229794	391409	32.77%	5.451%
19	9.334	1057	1062	1068	rVB	11132	15955	1.34%	0.222%
20	9.393	1068	1072	1077	rBV	9846	14342	1.20%	0.200%
21	9.752	1129	1133	1145	rVB5	7943	16000	1.34%	0.223%
22	9.905	1153	1159	1167	rVB2	15320	25681	2.15%	0.358%
23	10.504	1250	1261	1266	rBV	134748	240222	20.11%	3.346%
24	10.552	1266	1269	1276	rVB	35494	55020	4.61%	0.766%
25	12.175	1538	1545	1550	rBV	21651	34103	2.86%	0.475%
26	12.393	1576	1582	1592	rVB2	12413	21399	1.79%	0.298%
27	12.981	1673	1682	1692	rBV	582527	839233	70.27%	11.688%
28	13.828	1820	1826	1834	rBV	17669	28618	2.40%	0.399%
29	14.369	1911	1918	1924	rBV	208008	305414	25.57%	4.253%
30	15.863	2163	2172	2180	rBV	479145	690454	57.81%	9.616%
31	17.128	2380	2387	2393	rBV	268366	365452	30.60%	5.090%
32	18.022	2534	2539	2546	rBV	86762	125122	10.48%	1.743%
33	19.263	2746	2750	2760	rBV2	32313	53570	4.49%	0.746%
34	19.698	2819	2824	2830	rBV	998454	1194304	100.00%	16.633%
35	20.933	3031	3034	3040	rVB2	114747	154118	12.90%	2.146%
36	21.222	3079	3083	3088	rBV	305191	375196	31.42%	5.225%

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ClientSampleId :
B7-10-12

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-BP032321.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	23.510	3466	3472	3482	rVB2	201894	398508	33.37%	5.550%
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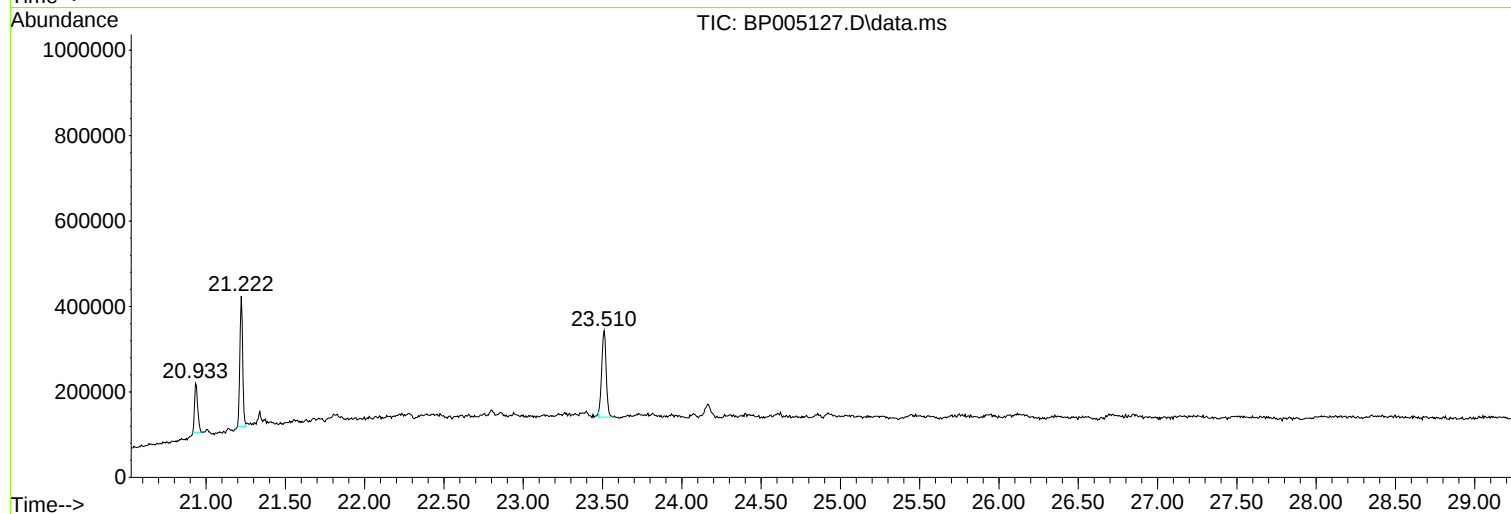
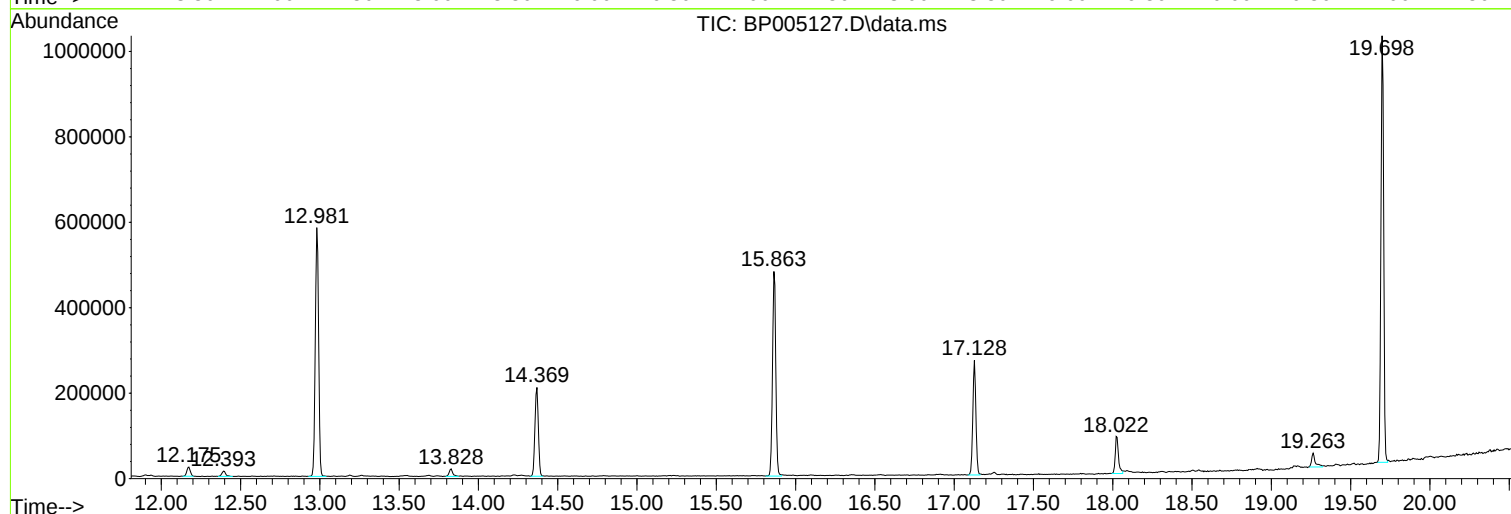
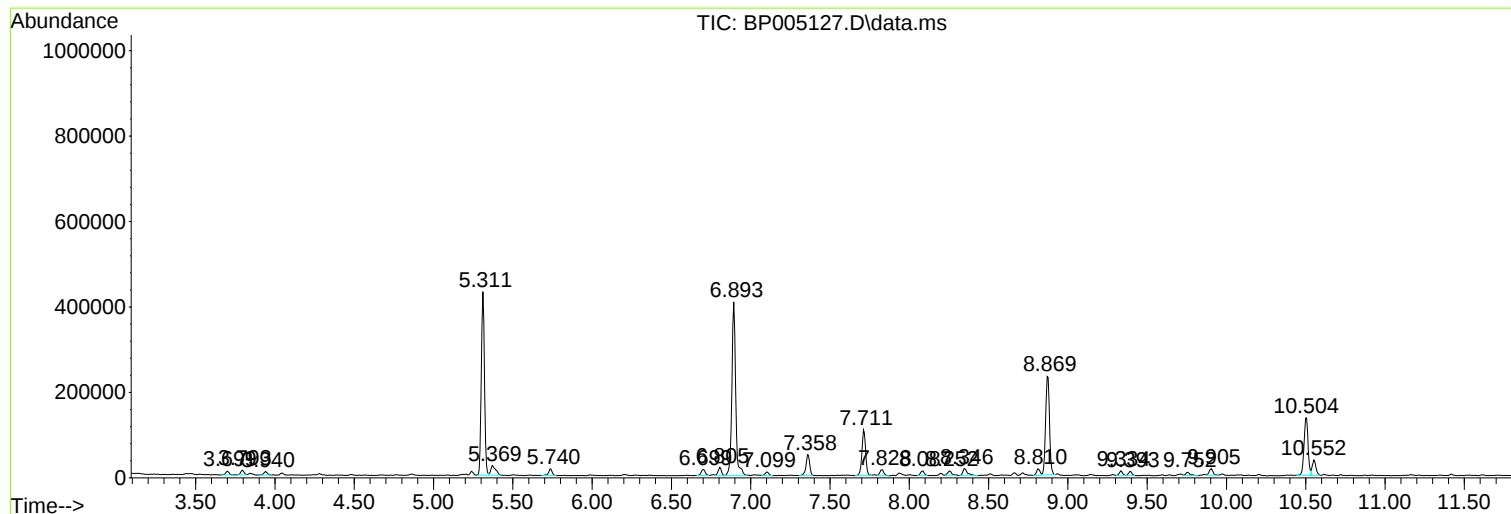
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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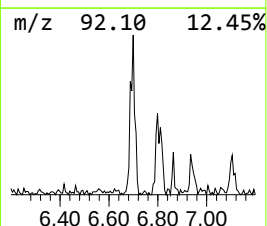
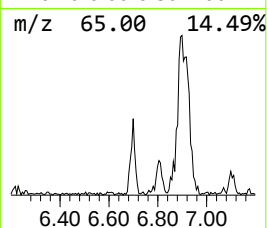
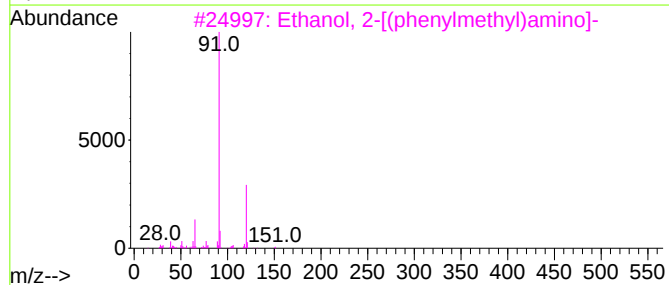
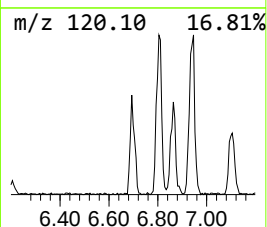
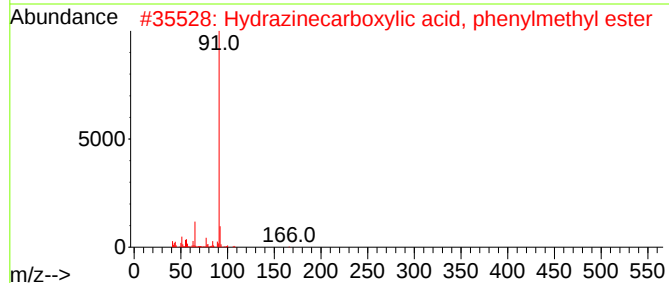
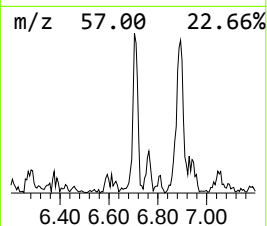
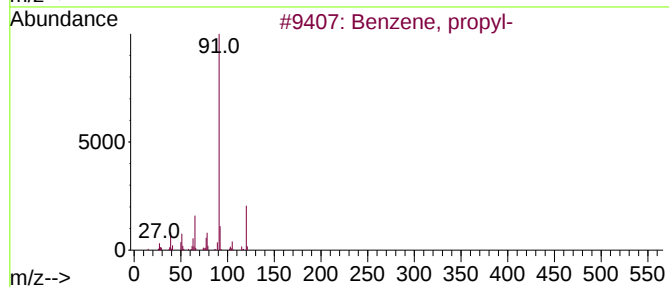
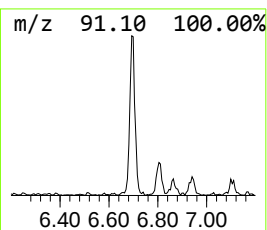
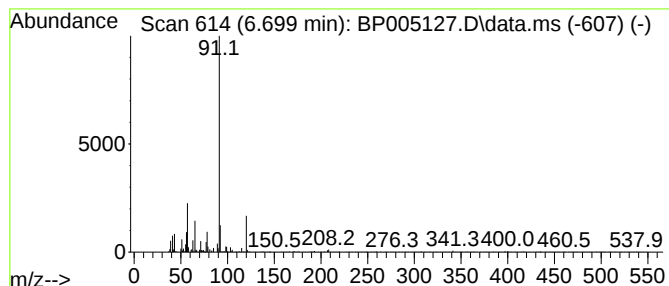
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.699	2.91 ng	25692	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	68
2			Hydrazinecarboxylic acid, phenyl...	166	C8H10N2O2	005331-43-1	49
3			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	47
4			N-Carboxybenzyl-S-benzylcysteine	345	C18H19NO4S	005619-12-5	46
5			Benzyl 2-p-toluenesulfonyl-.beta...	394	C19H22O7S	007252-07-5	46



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Instrument :
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ClientSampleId :
B7-10-12

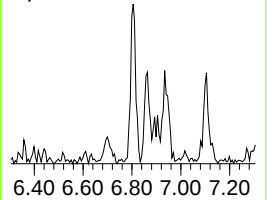
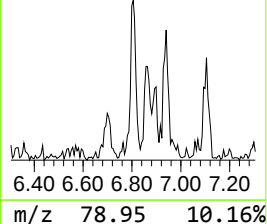
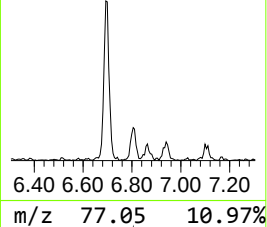
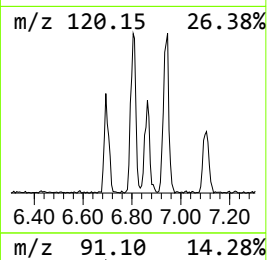
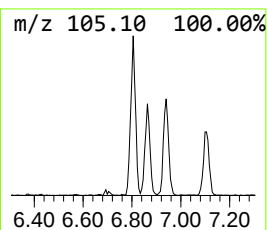
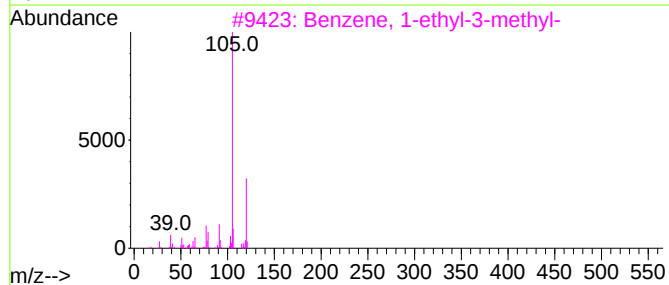
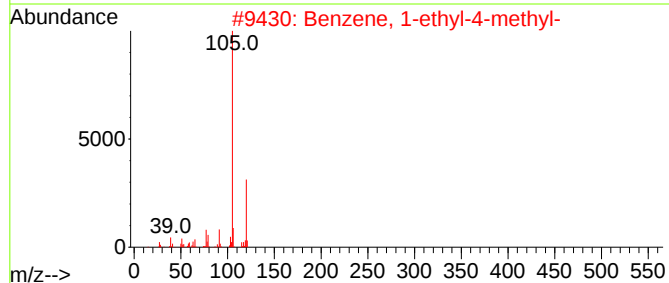
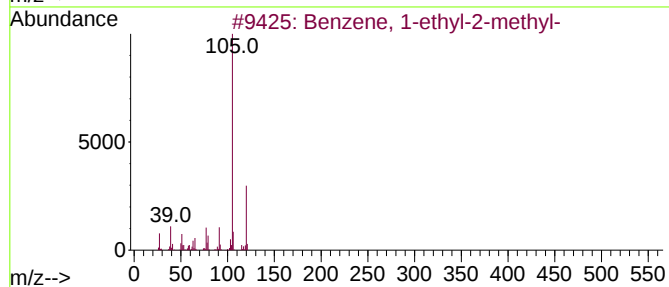
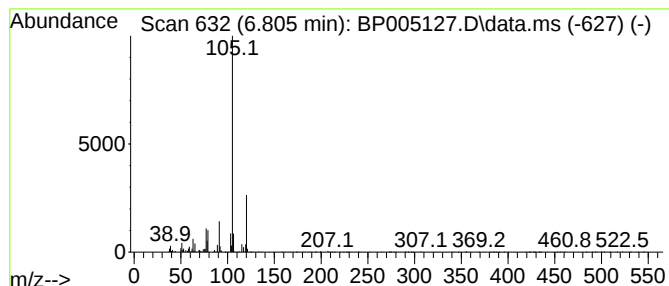
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.805	3.13 ng	27571	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	83



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ClientSampleId :
B7-10-12

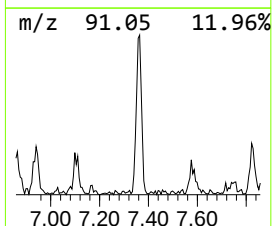
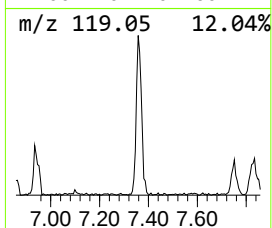
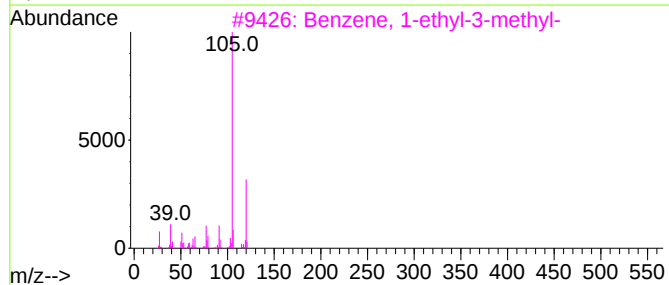
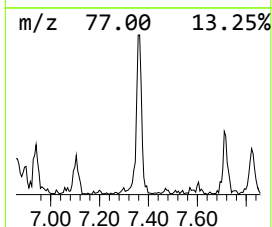
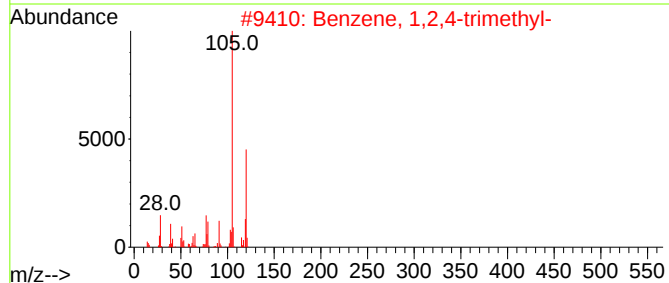
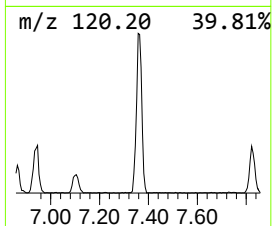
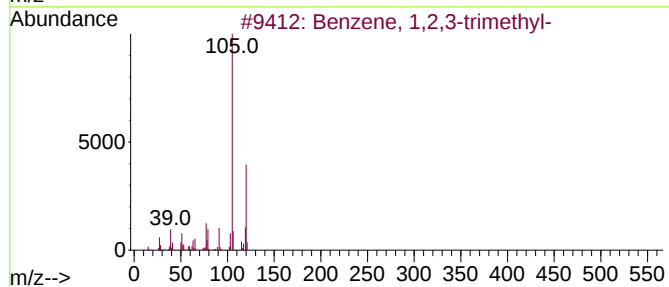
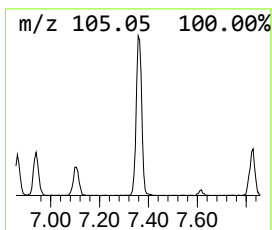
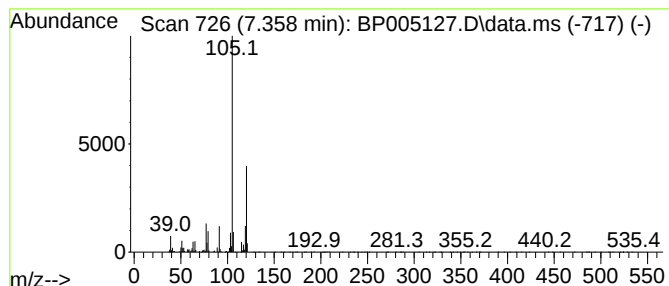
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.358	9.11 ng	80321	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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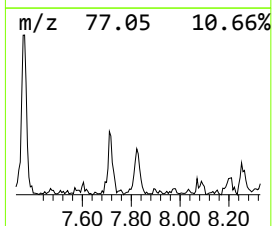
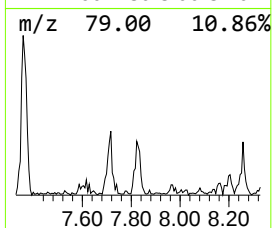
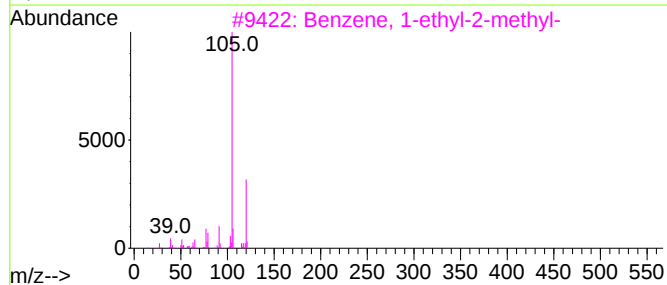
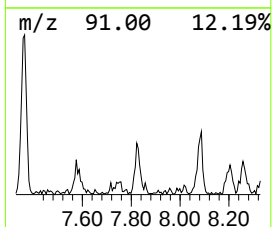
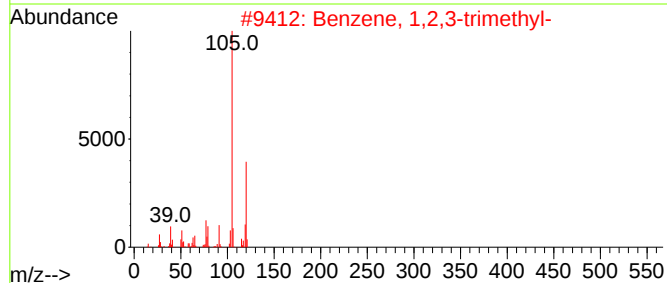
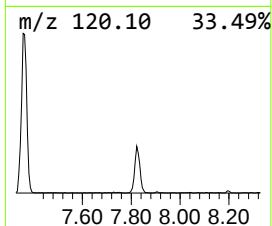
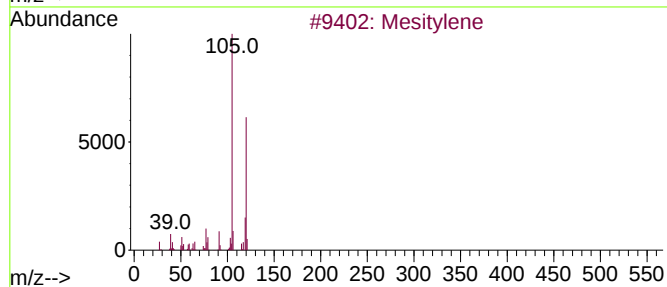
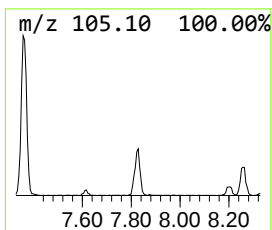
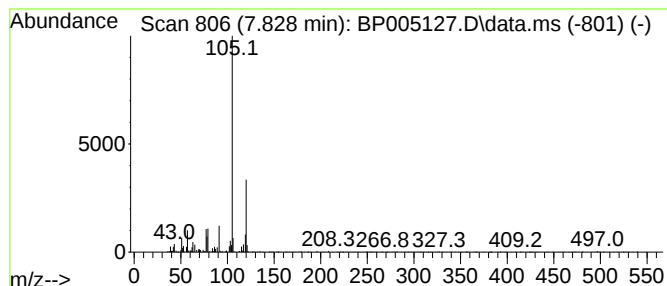
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Mesitylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	2.75 ng	24205	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	87
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	81
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	74



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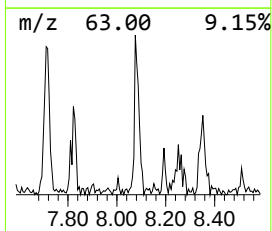
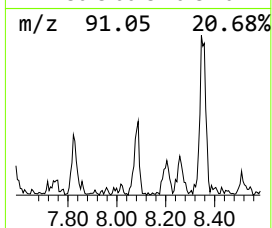
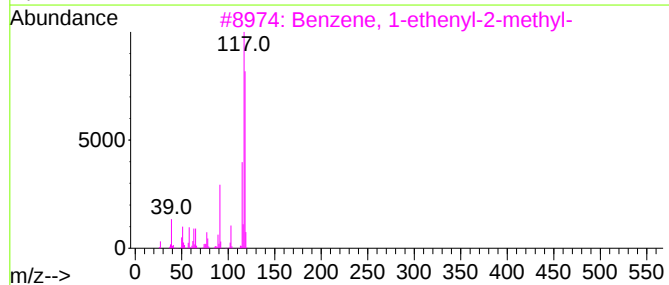
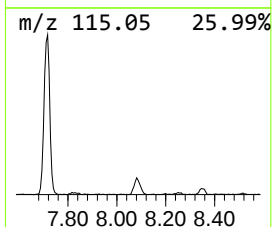
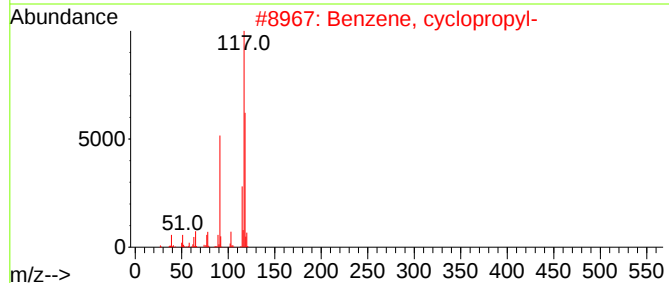
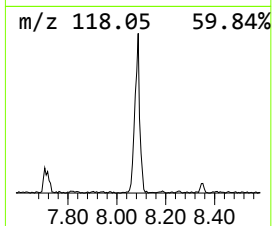
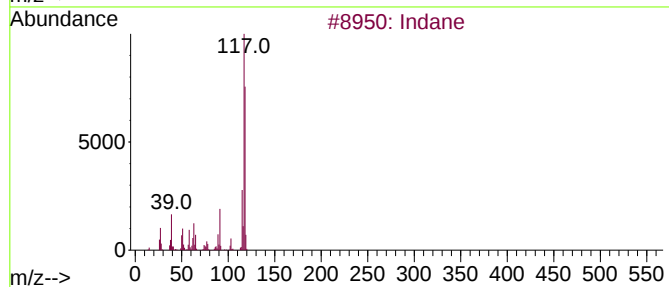
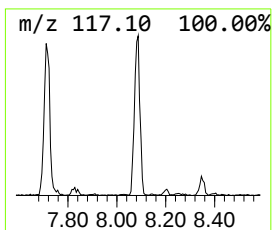
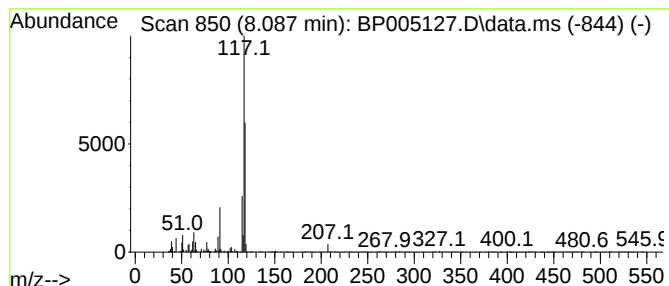
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Indane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.087	2.03 ng	17932	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	58
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	58
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58
4			Benzene, 1-(chloromethyl)-4-ethe...	152	C9H9Cl	001592-20-7	53
5			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
Acq On : 31 Mar 2021 15:37
Operator : CG/JU
Sample : M1836-16
Misc :
ALS Vial : 12 Sample Multiplier: 1

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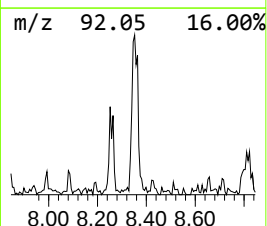
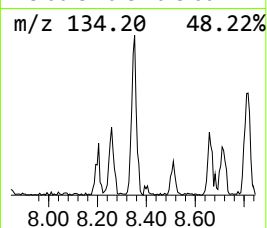
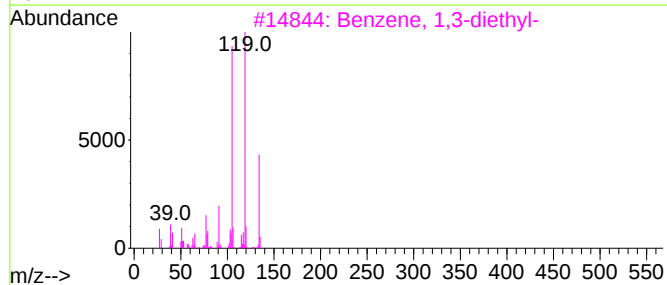
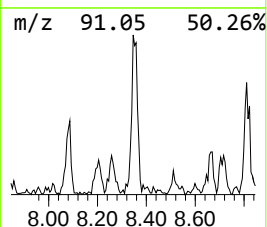
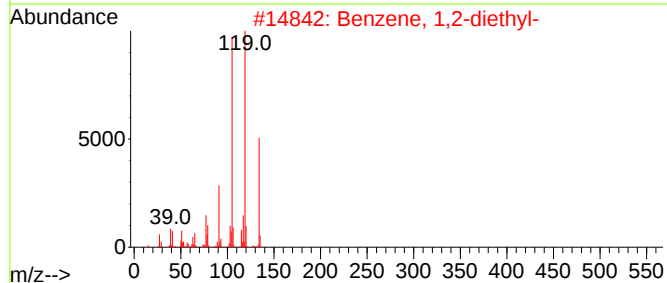
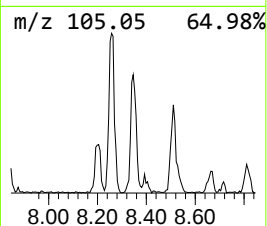
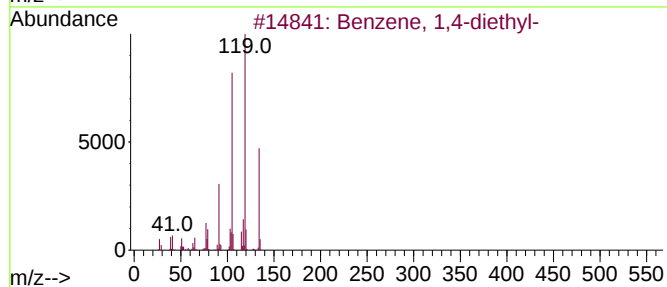
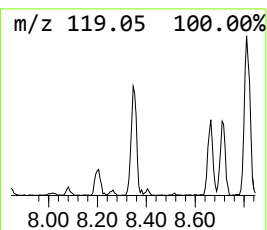
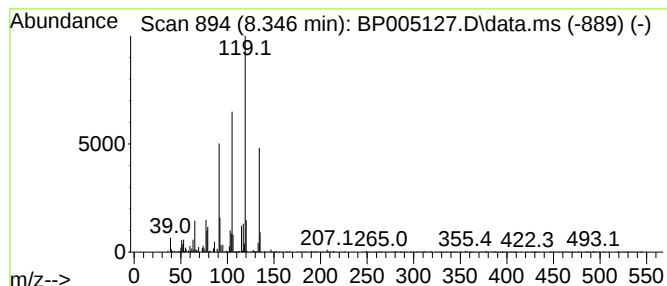
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,4-diethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.346	3.73 ng	32872	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
5			1-Methyl-1-silabenzocyclobutene	134	C8H10Si	160841-06-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
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Misc :
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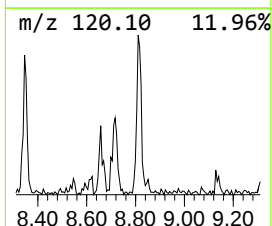
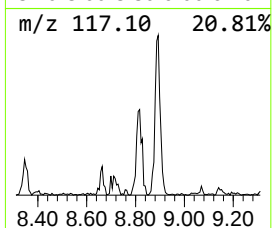
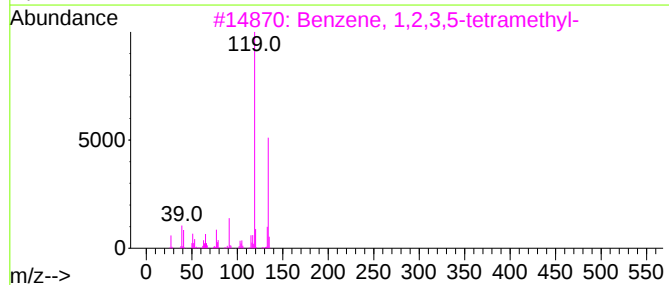
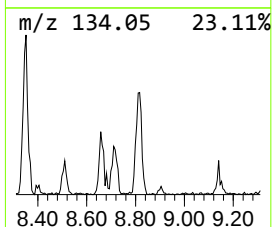
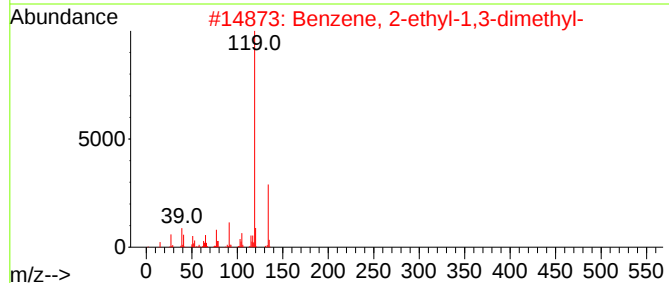
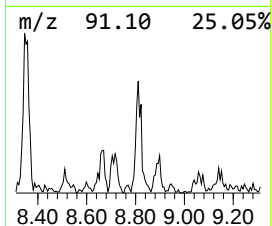
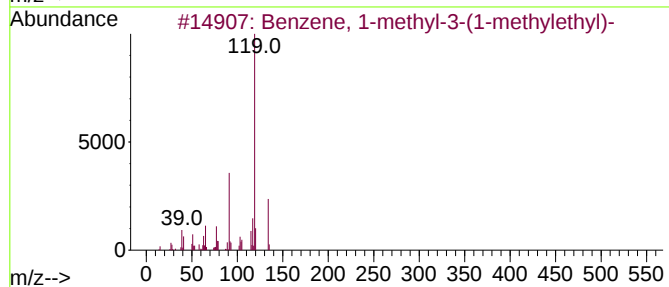
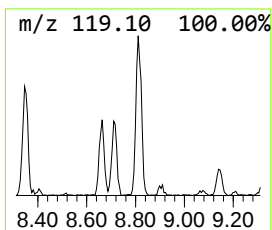
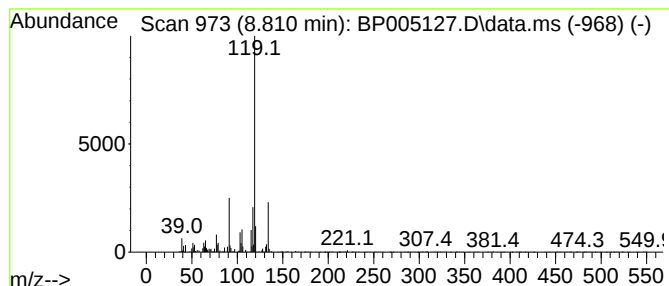
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-methyl-3-(1-meth... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.810	2.71 ng	23892	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
Acq On : 31 Mar 2021 15:37
Operator : CG/JU
Sample : M1836-16
Misc :
ALS Vial : 12 Sample Multiplier: 1

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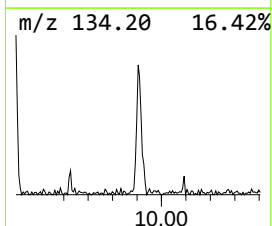
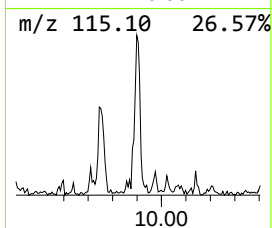
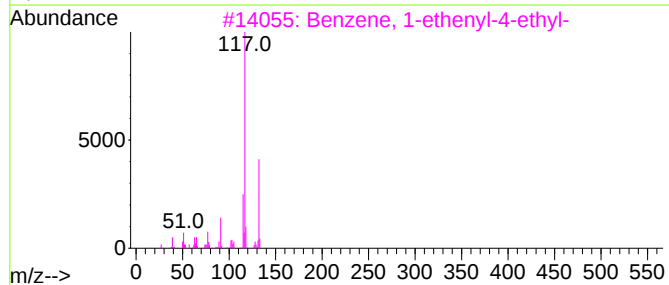
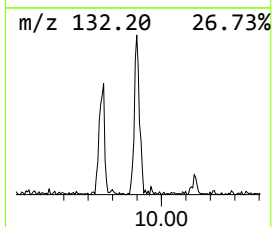
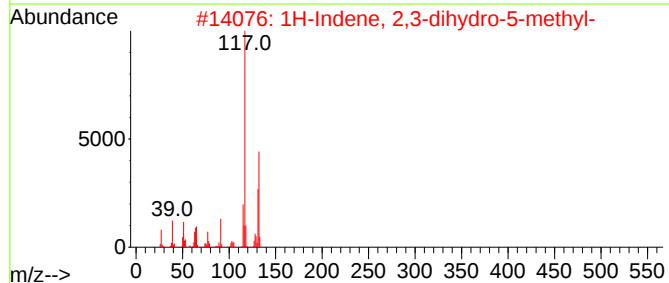
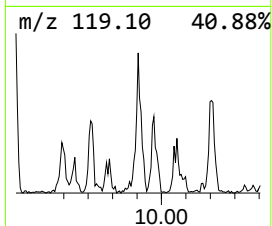
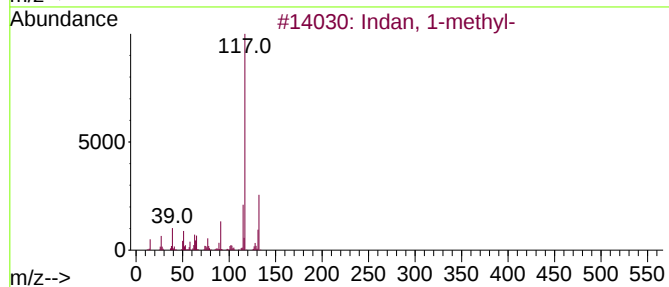
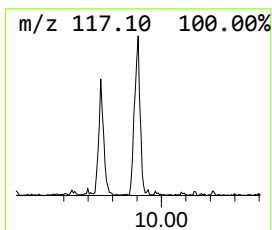
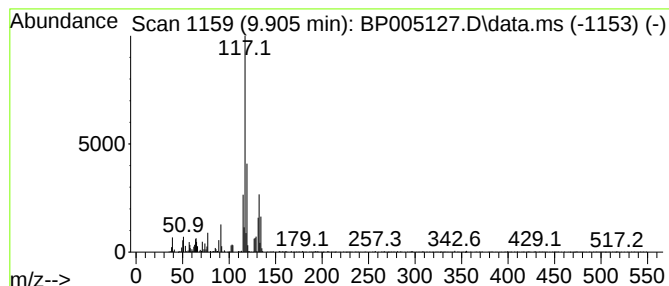
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Indan, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.905	2.14 ng	25681	Naphthalene-d8	10.505

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	70
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	58
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
Acq On : 31 Mar 2021 15:37
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Misc :
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Instrument :
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ClientSampleId :
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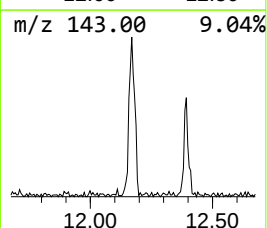
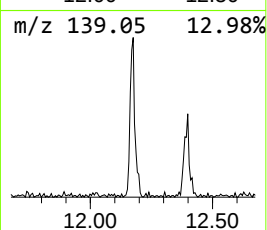
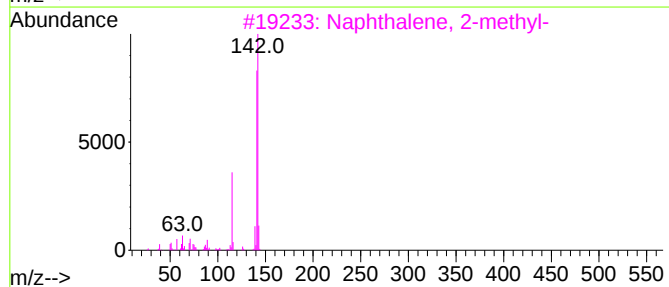
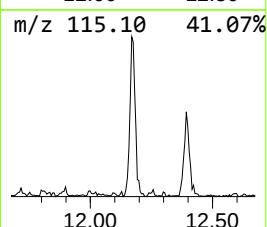
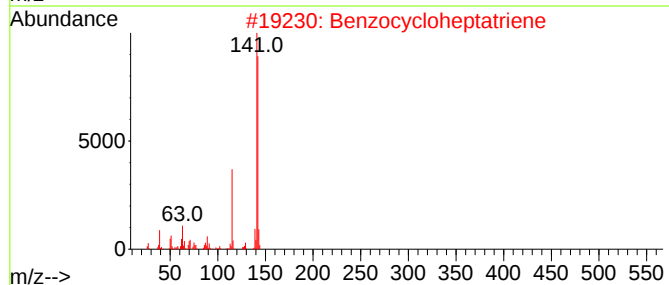
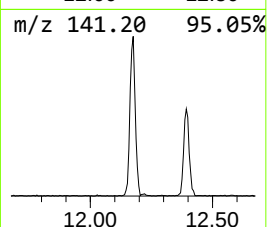
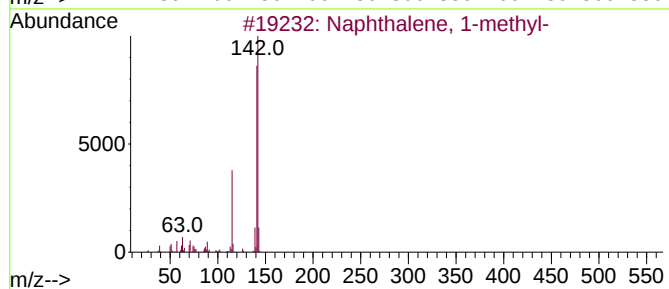
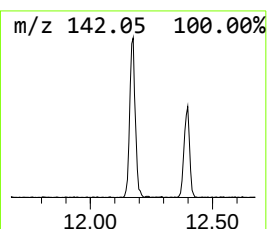
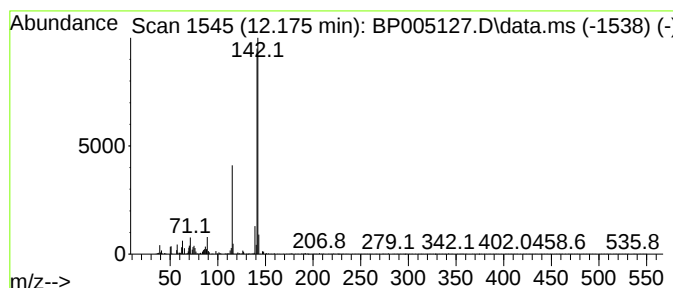
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.175	2.84 ng	34103	Naphthalene-d8	10.505

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2		Benzocycloheptatriene	142	C11H10	000264-09-5	91
3		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
Acq On : 31 Mar 2021 15:37
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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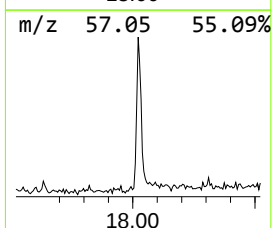
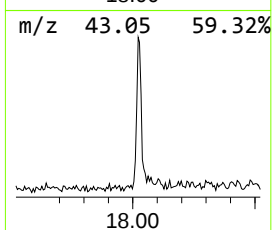
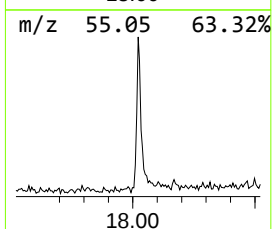
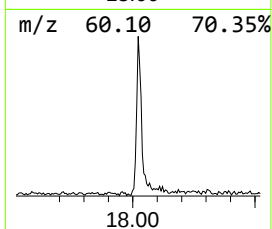
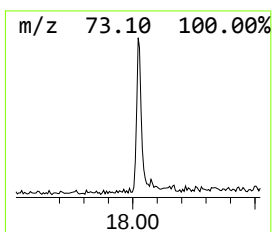
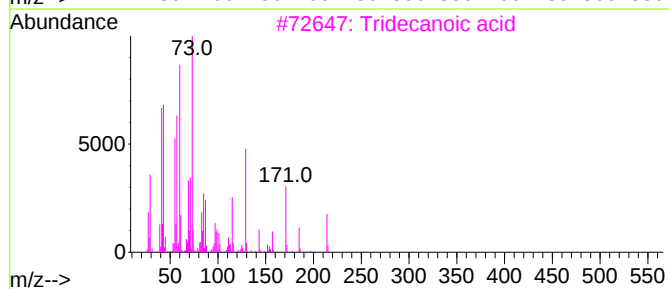
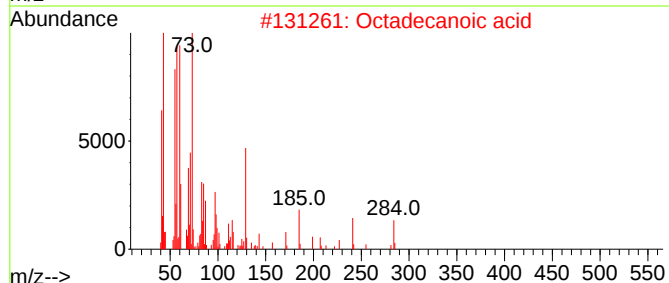
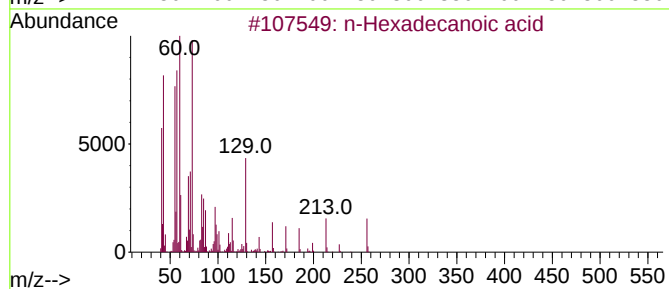
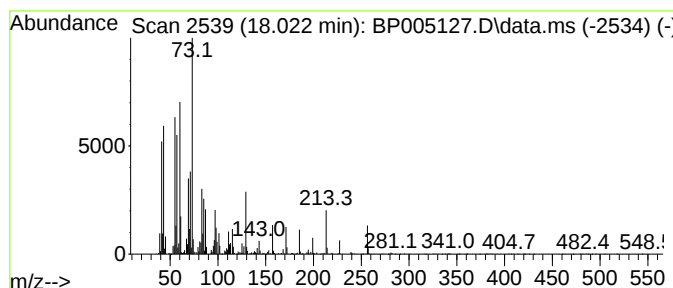
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	6.85 ng	125122	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	76
3			Tridecanoic acid	214	C13H26O2	000638-53-9	58
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	58
5			Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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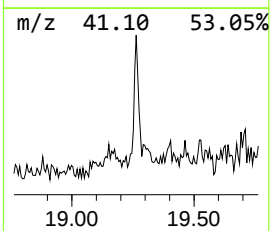
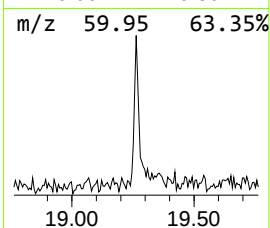
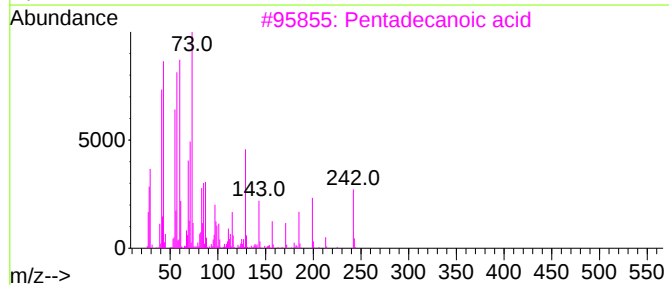
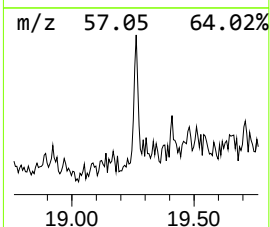
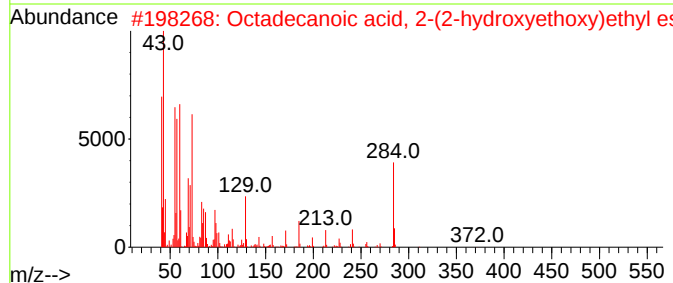
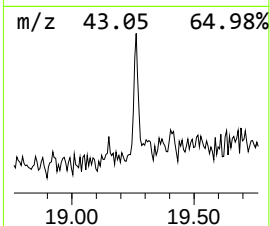
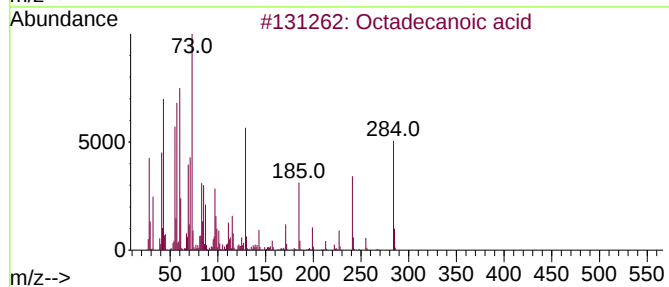
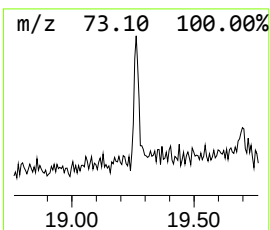
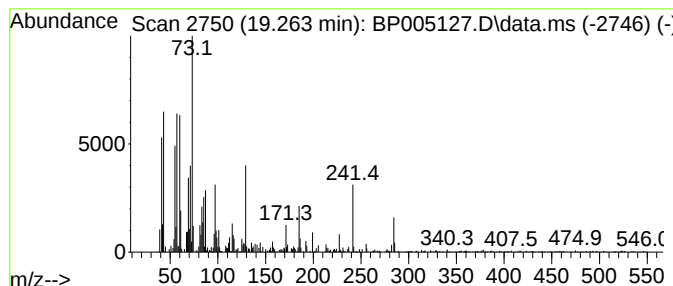
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.263	2.86 ng	53570	Chrysene-d12	21.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	93
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4			Tridecanoic acid	214	C13H26O2	000638-53-9	64
5			Eicosane	282	C20H42	000112-95-8	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
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Operator : CG/JU
Sample : M1836-16
Misc :
ALS Vial : 12 Sample Multiplier: 1

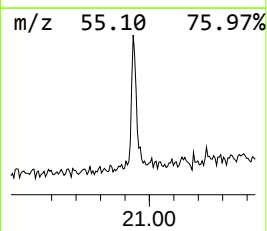
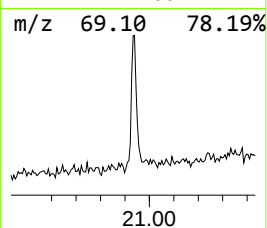
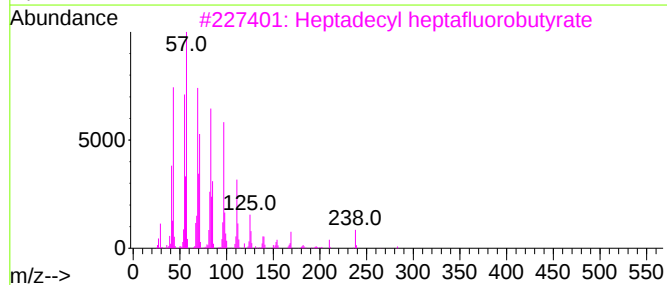
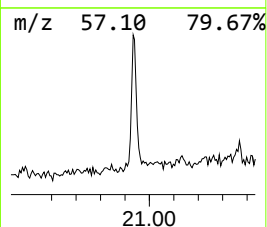
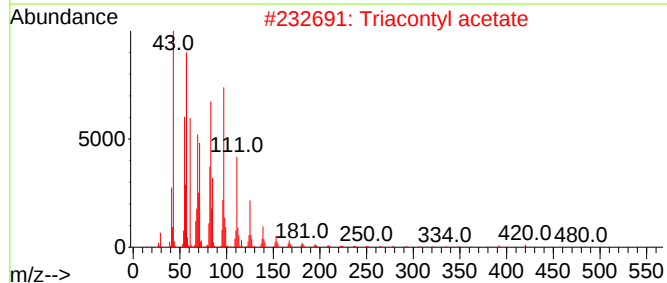
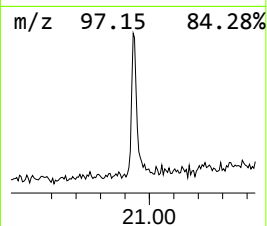
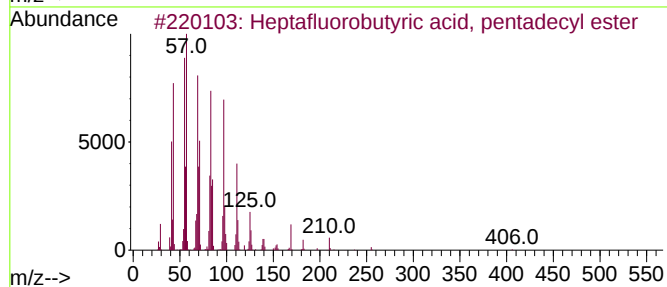
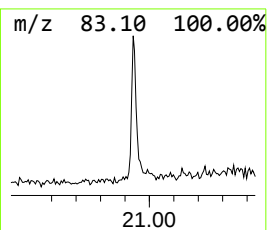
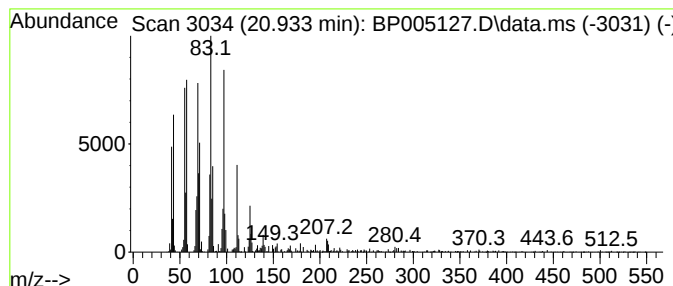
Instrument :
BNA_P
ClientSampleId :
B7-10-12

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptafluorobutyric acid, pe... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.933	8.22 ng	154118	Chrysene-d12	21.222	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	87
2	Triacontyl acetate	480	C32H64O2	041755-58-2	83
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	83
4	1-Docosene	308	C22H44	001599-67-3	68
5	5-Eicosene, (E)-	280	C20H40	074685-30-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
Data File : BP005127.D
Acq On : 31 Mar 2021 15:37
Operator : CG/JU
Sample : M1836-16
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
B7-10-12

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.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
Benzene, propyl-	6.699	2.9	ng	25692	1	7.710	176287	20.0
Benzene, 1-ethy...	6.805	3.1	ng	27571	1	7.710	176287	20.0
Benzene, 1,2,3-...	7.358	9.1	ng	80321	1	7.710	176287	20.0
Mesitylene	7.828	2.8	ng	24205	1	7.710	176287	20.0
Indane	8.087	2.0	ng	17932	1	7.710	176287	20.0
Benzene, 1,4-di...	8.346	3.7	ng	32872	1	7.710	176287	20.0
Benzene, 1-meth...	8.810	2.7	ng	23892	1	7.710	176287	20.0
Indan, 1-methyl-	9.905	2.1	ng	25681	2	10.505	240222	20.0
Naphthalene, 1-...	12.175	2.8	ng	34103	2	10.505	240222	20.0
n-Hexadecanoic ...	18.022	6.8	ng	125122	4	17.128	365452	20.0
Octadecanoic acid	19.263	2.9	ng	53570	5	21.222	375196	20.0
Heptafluorobuty...	20.933	8.2	ng	154118	5	21.222	375196	20.0