

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
 Data File : BF131404.D
 Acq On : 25 Nov 2022 19:47
 Operator : CG\JU
 Sample : N5741-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-31

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_F\Methods\8270-BF112122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131404.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.251	21	28	34	rBV	1347955	1730448	69.09%	9.406%
2	3.351	208	215	220	rBV2	20216	37424	1.49%	0.203%
3	4.492	403	409	415	rBV2	26093	37794	1.51%	0.205%
4	4.575	418	423	427	rBV	33873	35448	1.42%	0.193%
5	5.169	520	524	530	rVB	166999	171241	6.84%	0.931%
6	5.439	566	570	574	rVB	123891	107339	4.29%	0.583%
7	5.557	586	590	594	rBV	1214671	1130217	45.12%	6.143%
8	5.804	628	632	635	rBV	41920	37100	1.48%	0.202%
9	6.122	683	686	689	rVB	48104	36915	1.47%	0.201%
10	6.410	732	735	739	rVB	34938	29390	1.17%	0.160%
11	6.563	756	761	764	rBV	878636	737790	29.46%	4.010%
12	6.639	770	774	777	rVB	58664	45882	1.83%	0.249%
13	6.775	794	797	802	rVB	35623	34598	1.38%	0.188%
14	6.957	824	828	831	rBV	567931	510993	20.40%	2.778%
15	7.010	834	837	840	rVB	55948	42225	1.69%	0.230%
16	7.133	855	858	861	rVB	196364	152412	6.08%	0.828%
17	7.522	919	924	927	rVB	1561796	1442647	57.60%	7.842%
18	7.804	968	972	976	rBV2	25007	26747	1.07%	0.145%
19	7.974	997	1001	1005	rBV2	61981	67019	2.68%	0.364%
20	8.245	1040	1047	1049	rBV	762062	679861	27.14%	3.695%
21	8.263	1049	1050	1053	rVB	78217	46903	1.87%	0.255%
22	8.463	1081	1084	1086	rBV	32291	28873	1.15%	0.157%
23	8.710	1122	1126	1131	rBV5	26794	44538	1.78%	0.242%
24	8.780	1136	1138	1140	rBV	26908	26695	1.07%	0.145%
25	9.057	1181	1185	1189	rBV3	121300	126440	5.05%	0.687%
26	9.133	1195	1198	1203	rBV3	29472	44521	1.78%	0.242%
27	9.186	1204	1207	1212	rVB6	34516	39861	1.59%	0.217%
28	9.245	1212	1217	1221	rBV8	18864	31652	1.26%	0.172%
29	9.327	1226	1231	1233	rVV	2720964	2504766	100.00%	13.615%
30	9.351	1233	1235	1241	rVB2	52743	51233	2.05%	0.278%
31	9.492	1256	1259	1262	rVB	95027	87370	3.49%	0.475%
32	9.586	1270	1275	1281	rBV8	43617	85927	3.43%	0.467%
33	9.674	1287	1290	1293	rVB5	38319	48337	1.93%	0.263%
34	9.751	1296	1303	1306	rBV9	20902	44521	1.78%	0.242%
35	9.857	1318	1321	1325	rBV3	51394	55791	2.23%	0.303%
36	9.904	1325	1329	1334	rVB	177002	176831	7.06%	0.961%

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37	10.010	1343	1347	1350	rBV	911346	740474	29.56%	4.025%
38	10.204	1378	1380	1384	rVB5	33626	29473	1.18%	0.160%
39	10.321	1396	1400	1407	rBV	141882	312077	12.46%	1.696%
40	10.433	1417	1419	1423	rVB3	38277	33060	1.32%	0.180%
41	10.627	1449	1452	1456	rVB2	44433	61403	2.45%	0.334%
42	10.727	1465	1469	1472	rVB	117735	102078	4.08%	0.555%
43	10.763	1472	1475	1477	rBV3	27287	27688	1.11%	0.150%
44	10.804	1477	1482	1485	rBV	1938265	1675633	66.90%	9.108%
45	11.157	1540	1542	1545	rVB4	39365	32400	1.29%	0.176%
46	11.380	1578	1580	1584	rVB2	37170	38616	1.54%	0.210%
47	11.504	1598	1601	1605	rBV	773755	701776	28.02%	3.815%
48	12.027	1687	1690	1697	rVB	307666	276819	11.05%	1.505%
49	12.286	1732	1734	1737	rVB	54589	43593	1.74%	0.237%
50	12.815	1821	1824	1830	rVB3	73560	78967	3.15%	0.429%
51	13.104	1869	1873	1876	rBV	2571360	2344778	93.61%	12.745%
52	13.998	2022	2025	2037	rVB	165442	189508	7.57%	1.030%
53	14.162	2050	2053	2057	rVB	576137	481421	19.22%	2.617%
54	14.262	2067	2070	2075	rVB	90728	89272	3.56%	0.485%
55	15.321	2246	2250	2254	rBV3	21739	28057	1.12%	0.153%
56	15.698	2309	2314	2319	rBV	345217	451705	18.03%	2.455%
57	16.209	2397	2401	2405	rVB2	19520	28297	1.13%	0.154%
58	16.756	2490	2494	2499	rBV3	18309	31656	1.26%	0.172%
59	17.409	2600	2605	2611	rBV8	15706	29079	1.16%	0.158%
60	18.186	2730	2737	2745	rVB9	12338	31881	1.27%	0.173%

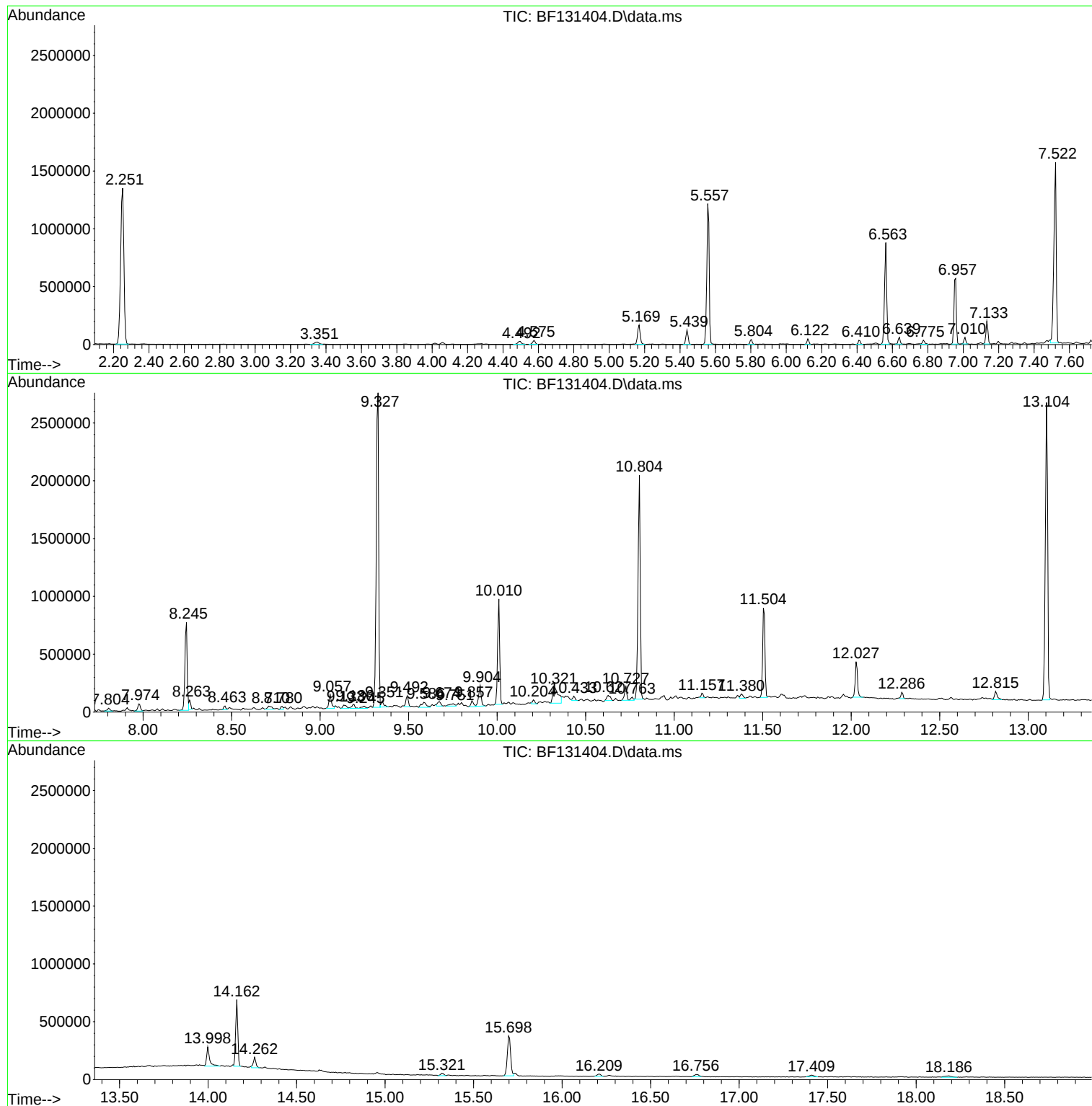
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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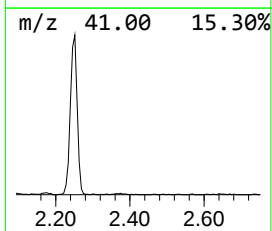
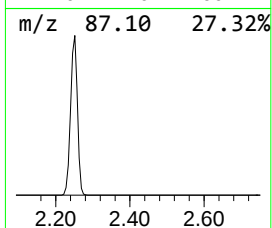
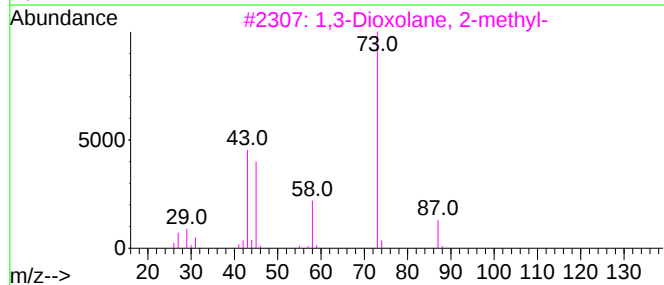
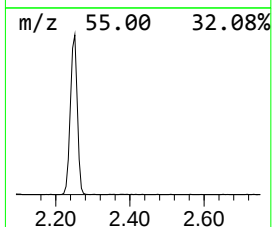
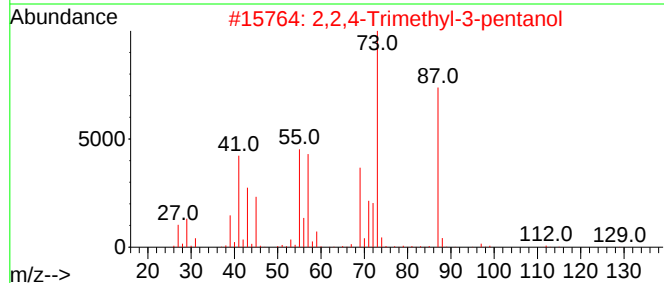
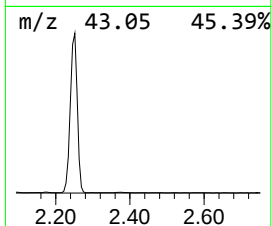
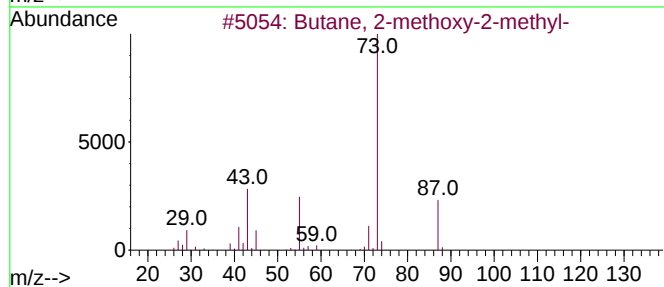
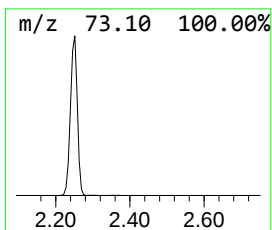
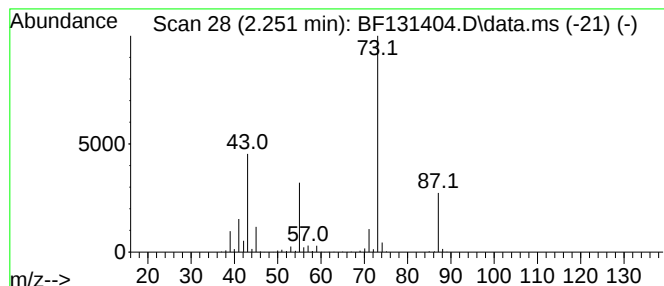
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.251	67.73 ng	1730450	1,4-Dichlorobenzene-d4	6.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	37
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	12



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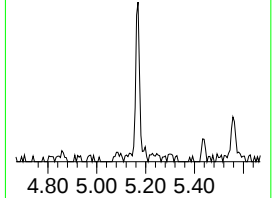
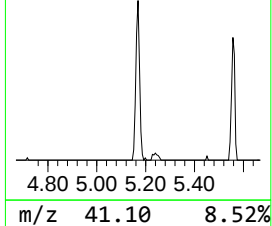
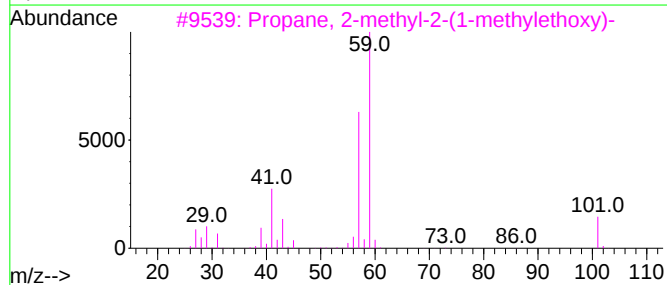
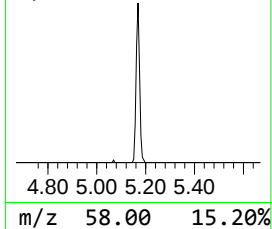
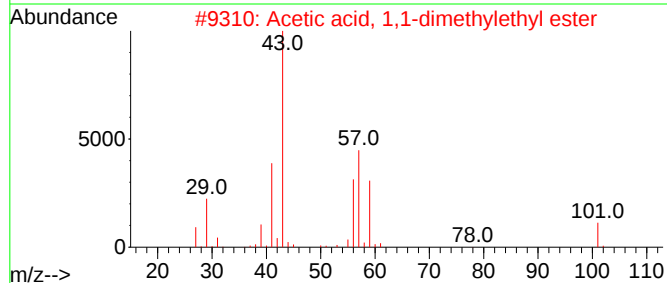
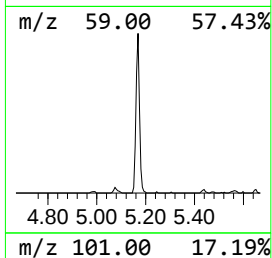
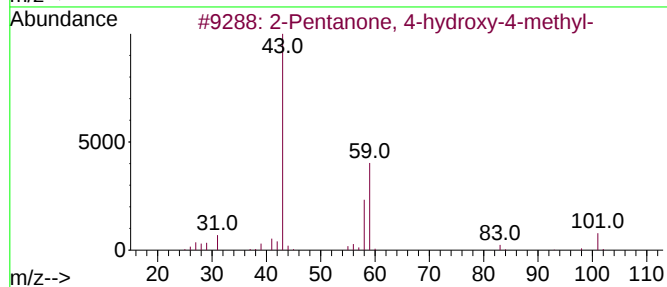
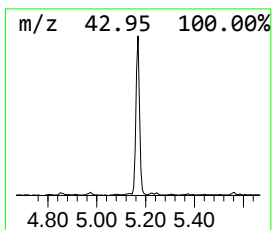
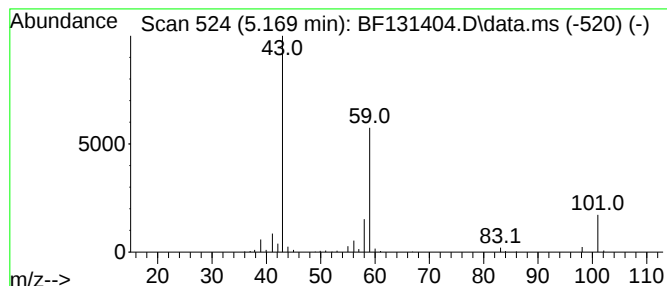
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.169	6.70 ng	171241	1,4-Dichlorobenzene-d4	6.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		



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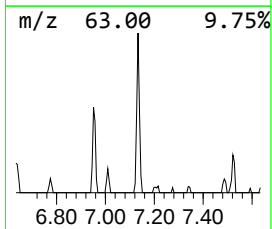
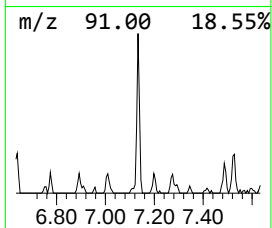
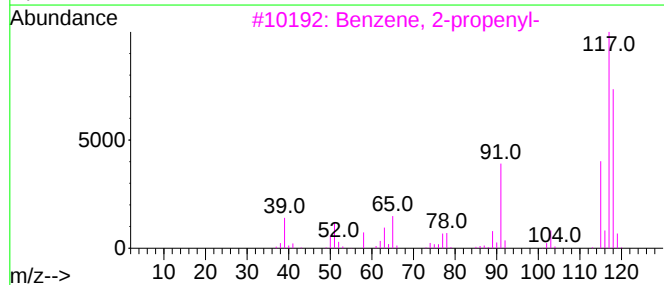
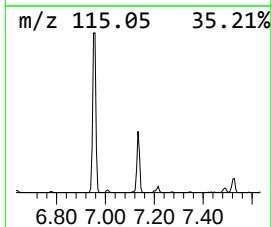
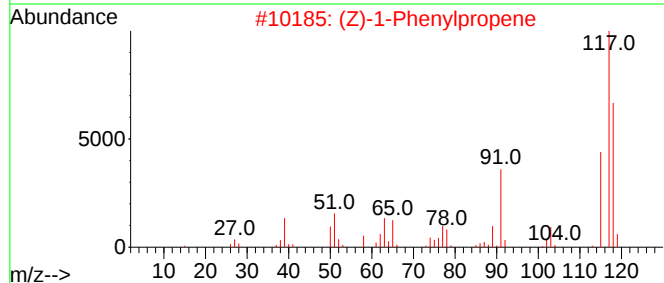
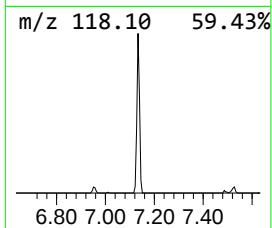
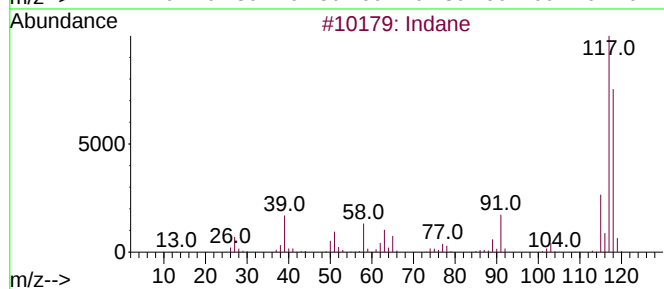
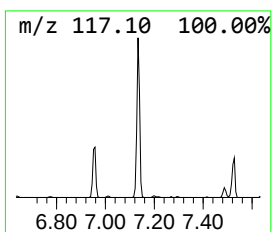
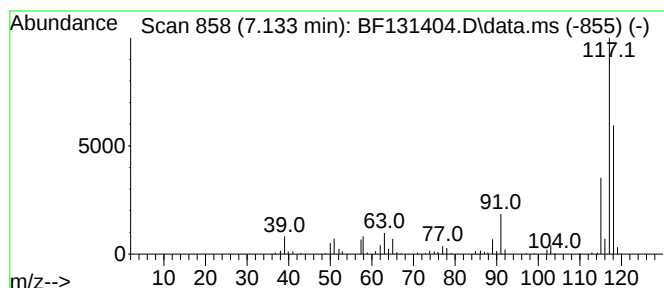
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.133	5.97 ng	152412	1,4-Dichlorobenzene-d4	6.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	91
2			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	74
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	72
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64



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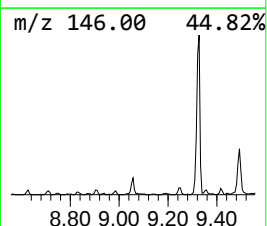
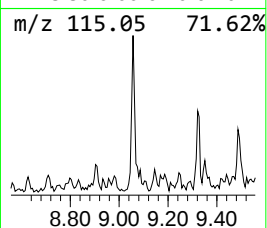
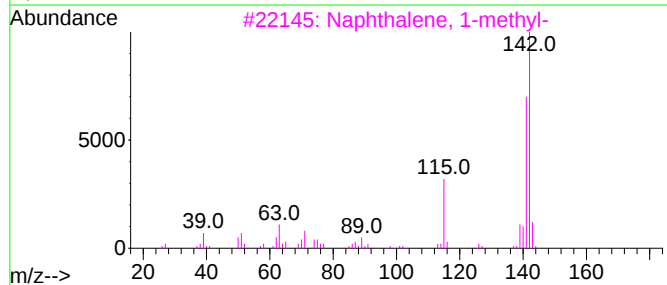
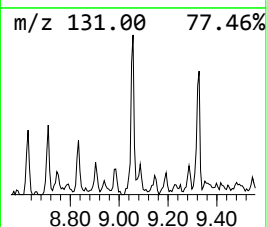
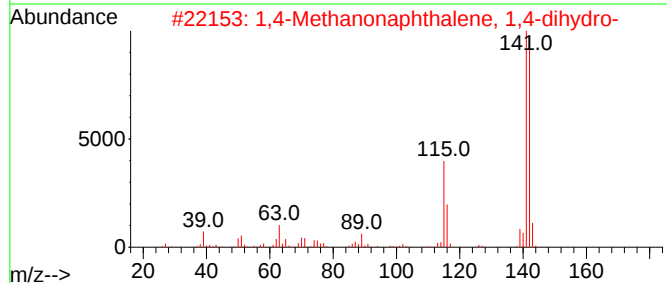
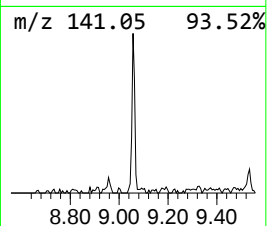
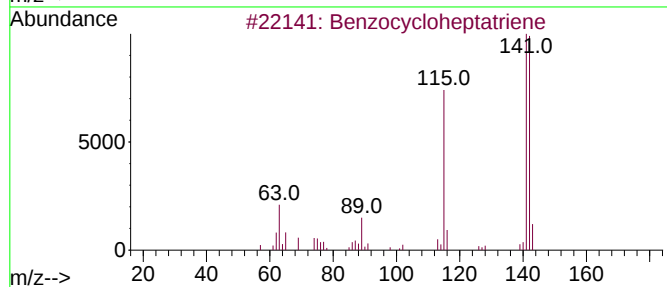
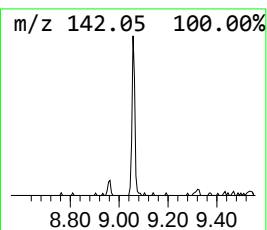
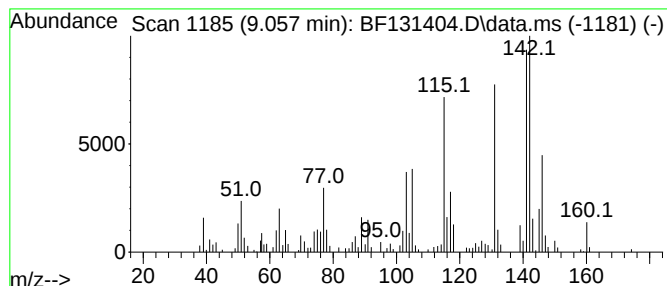
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzocycloheptatriene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.057	3.72 ng	126440	Naphthalene-d8	8.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzocycloheptatriene	142	C11H10	000264-09-5	70
2			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	55
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	55
4			1-Methylindan-2-one	146	C10H10O	035587-60-1	49
5			Benzeneacetonitrile, 4-cyano-	142	C9H6N2	000876-31-3	42



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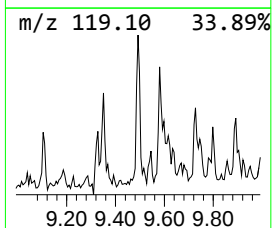
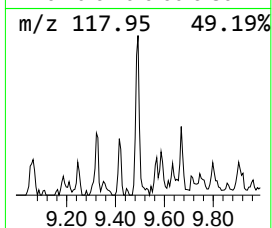
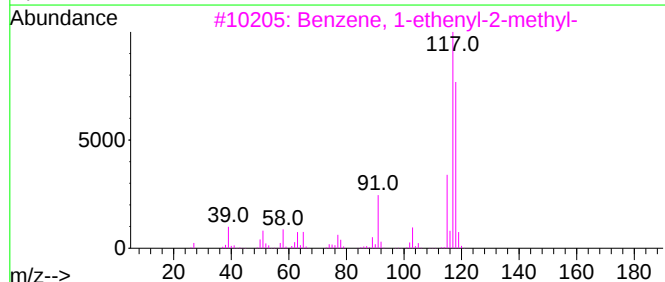
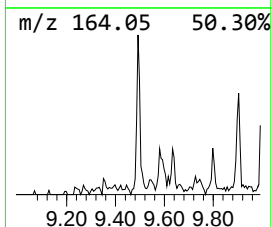
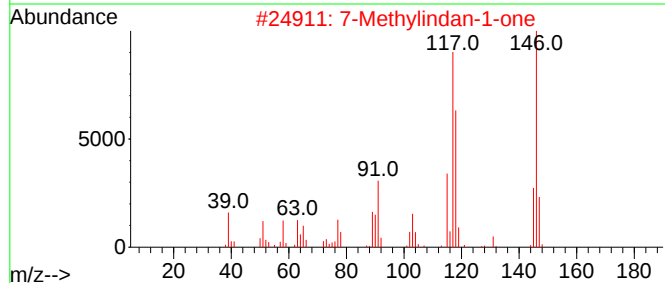
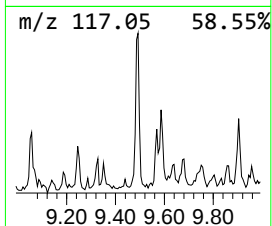
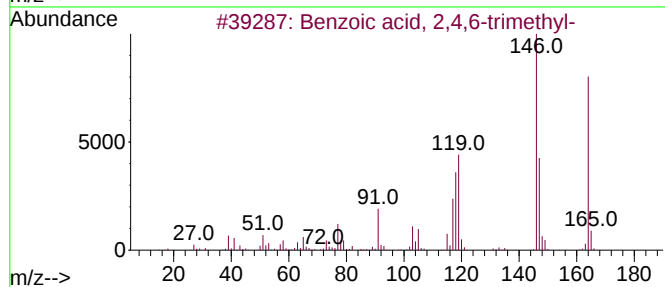
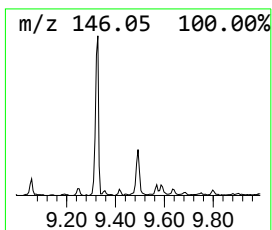
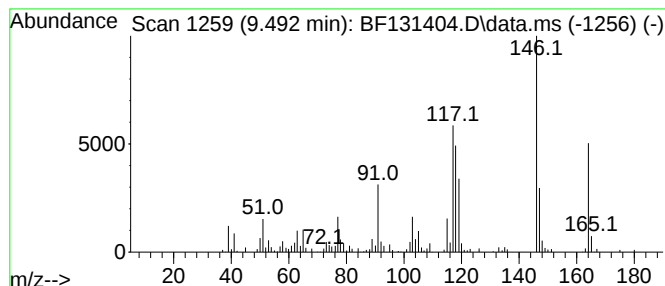
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.492	2.36 ng	87370	Acenaphthene-d10	10.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	64
2			7-Methylindan-1-one	146	C10H10O	039627-61-7	58
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	47
4			4-Fluorophthalonitrile	146	C8H3FN2	065610-14-2	38
5			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-31

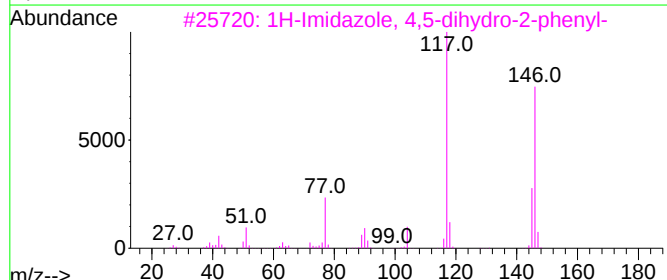
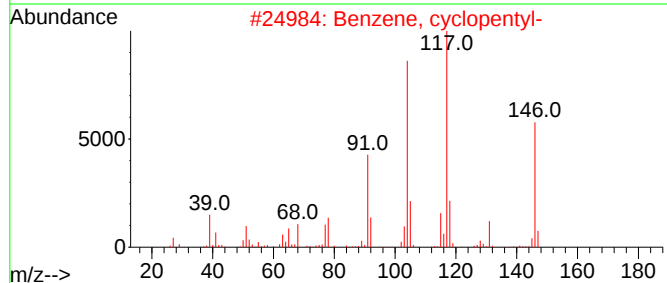
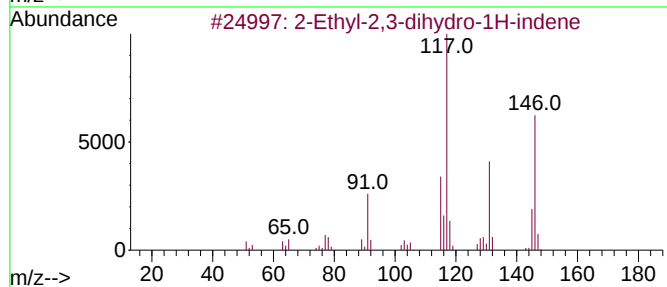
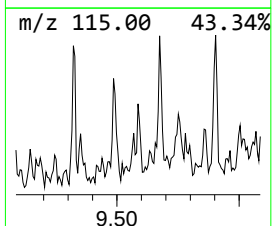
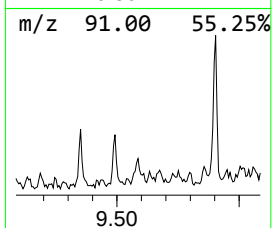
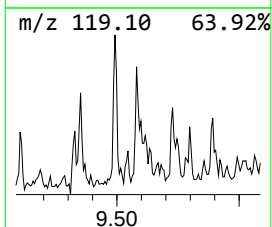
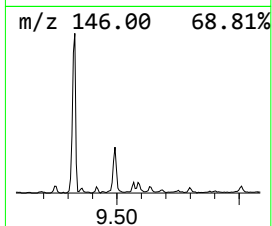
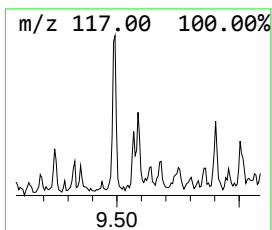
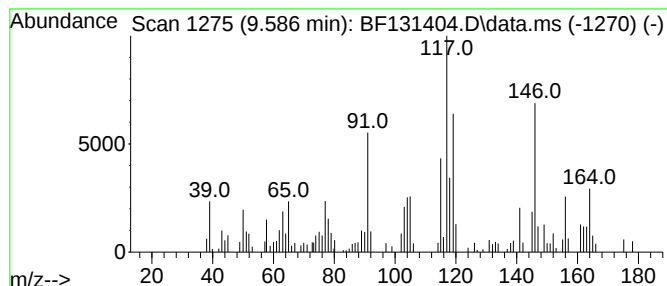
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.586 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.586	2.32 ng	85927	Acenaphthene-d10	10.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	43		
2	Benzene, cyclopentyl-	146	C11H14	000700-88-9	38		
3	1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	38		
4	m-Aminophenylacetylene	117	C8H7N	054060-30-9	38		
5	Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
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Sample : N5741-12
Misc :
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Instrument :
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ClientSampleId :
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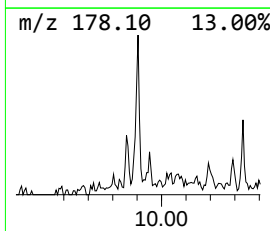
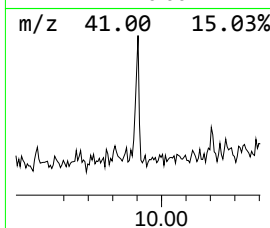
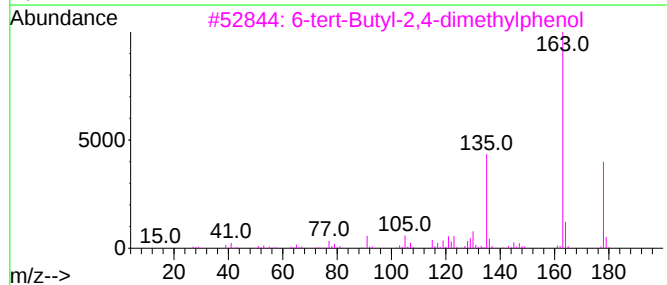
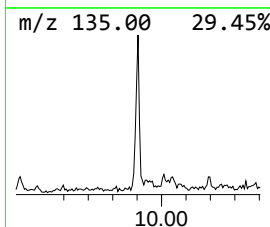
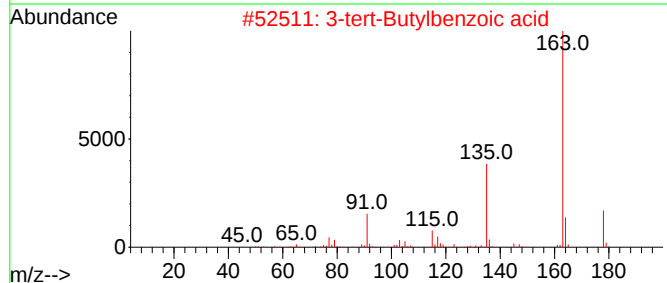
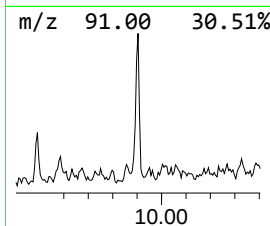
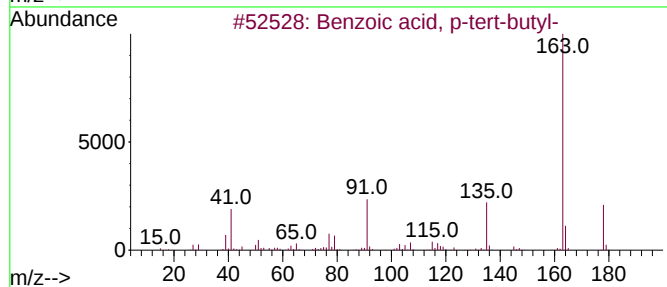
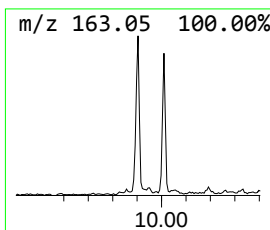
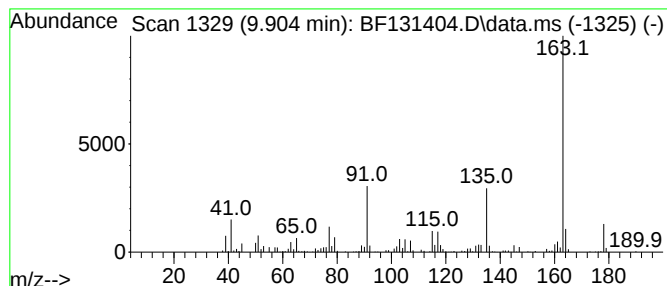
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzoic acid, p-tert-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.904	4.78 ng	176831	Acenaphthene-d10	10.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	90
2			3-tert-Butylbenzoic acid	178	C11H14O2	007498-54-6	80
3			6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	68
4			Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	68
5			Benzoic acid, 4-acetyl-, methyl ...	178	C10H10O3	003609-53-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

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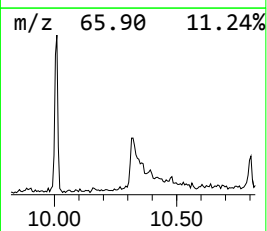
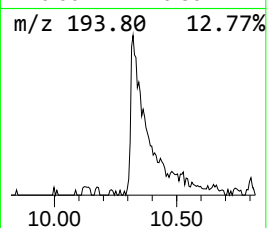
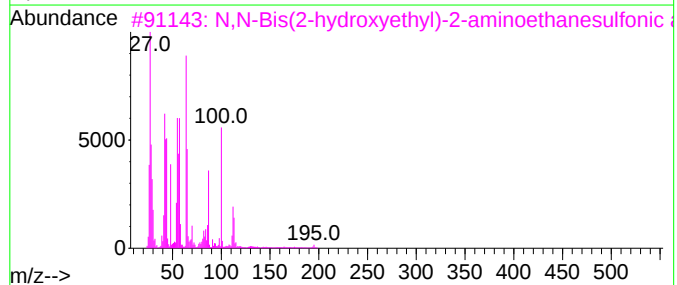
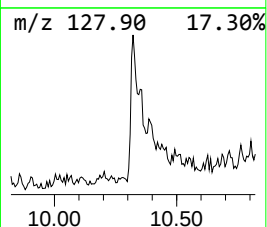
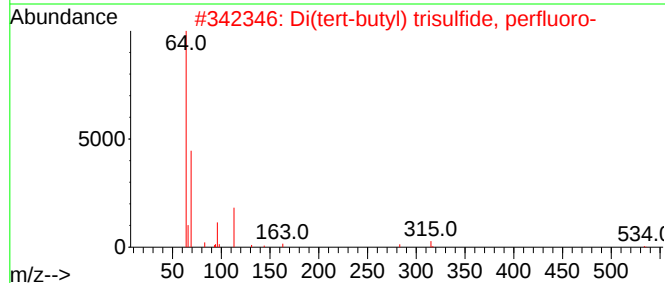
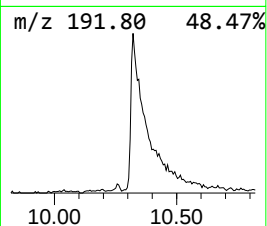
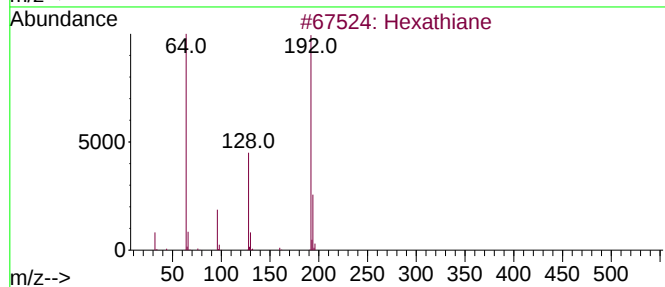
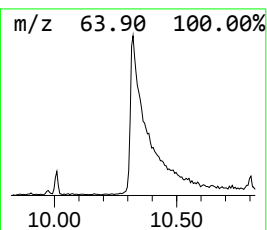
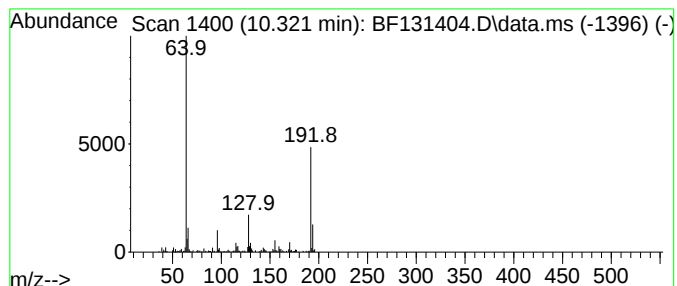
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexathiane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.321	8.43 ng	312077	Acenaphthene-d10	10.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexathiane	192	S6	013798-23-7	91
2			Di(tert-butyl) trisulfide, perfluoro-	534	C8F18S3	016005-64-4	39
3			N,N-Bis(2-hydroxyethyl)-2-aminoethanesulfonic acid	213	C6H15NO5S	010191-18-1	9
4			Tetrafluoroformamidine	116	CF4N2	014362-70-0	7
5			Sulfur dioxide	64	O2S	007446-09-5	4



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

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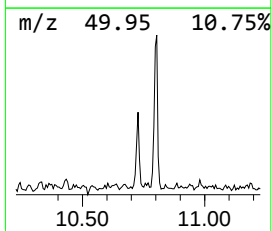
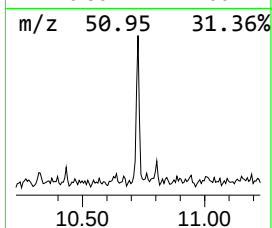
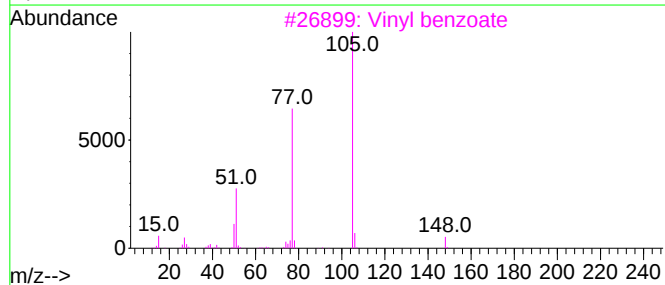
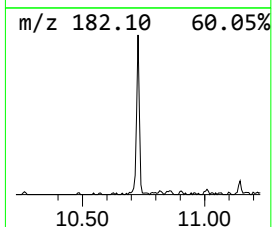
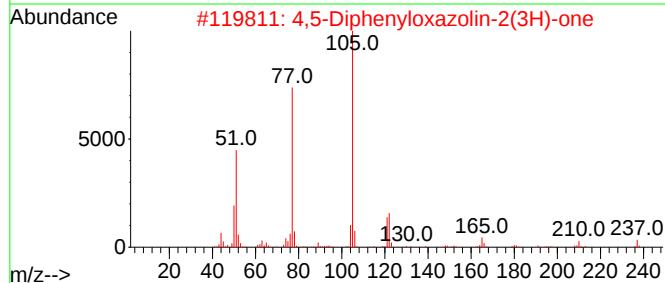
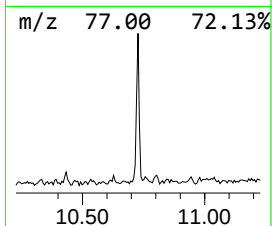
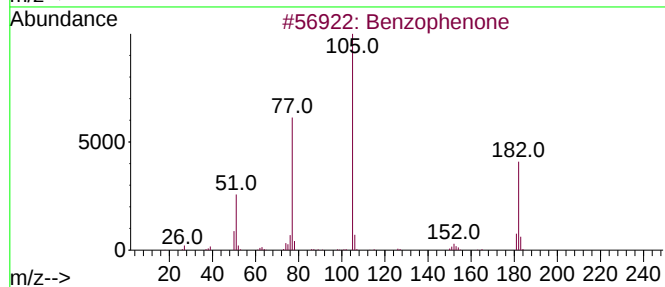
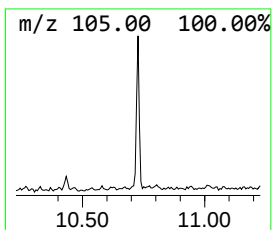
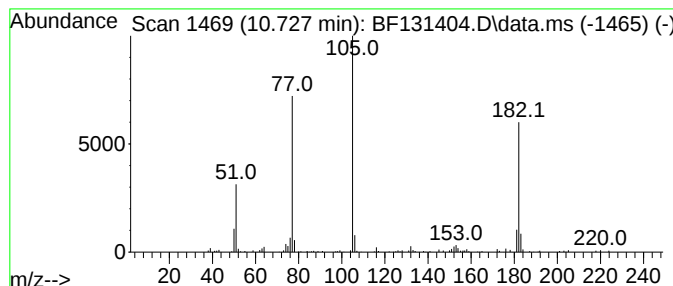
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.727	2.76 ng	102078	Acenaphthene-d10	10.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	95
2			4,5-Diphenyloxazolin-2(3H)-one	237	C15H11NO2	005014-83-5	47
3			Vinyl benzoate	148	C9H8O2	000769-78-8	47
4			Methyl 2-benzoyl-4-cyanobutanoate	231	C13H13NO3	1010486-83-8	47
5			2-Phenyl-4-ethylidene-2-oxazolin...	187	C11H9NO2	037791-41-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

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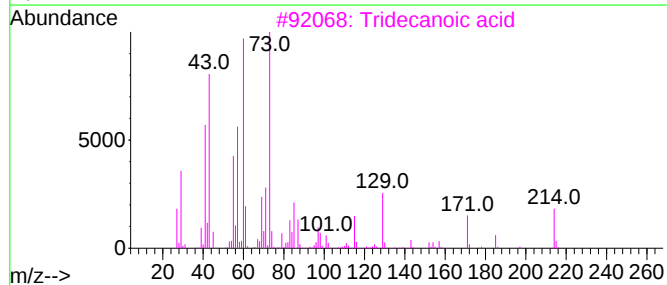
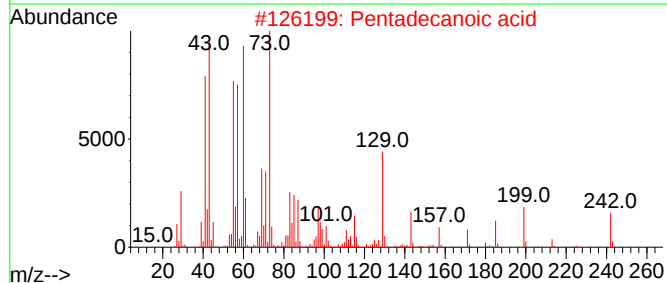
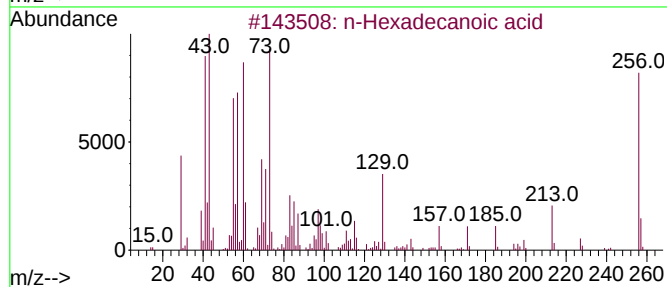
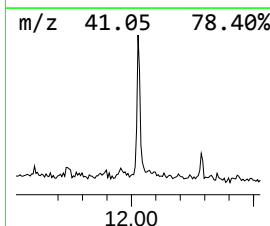
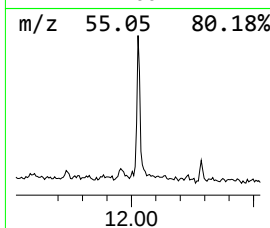
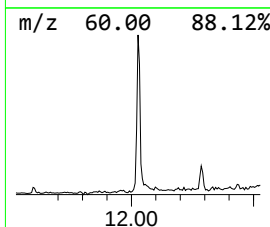
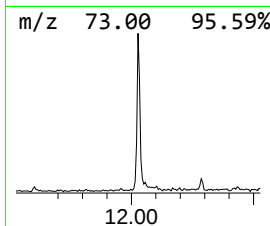
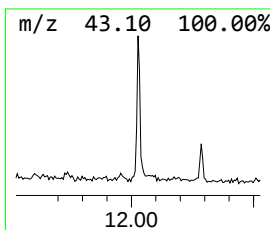
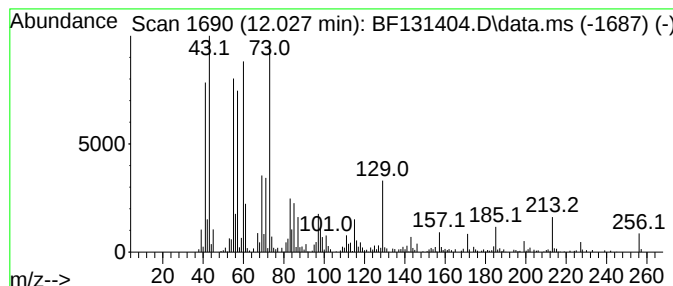
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.027	7.89 ng	276819	Phenanthrene-d10	11.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
5			Undecanoic acid	186	C11H22O2	000112-37-8	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
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Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

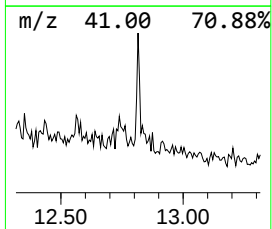
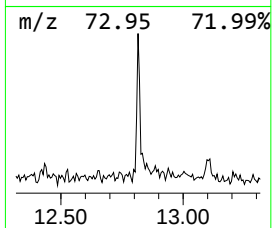
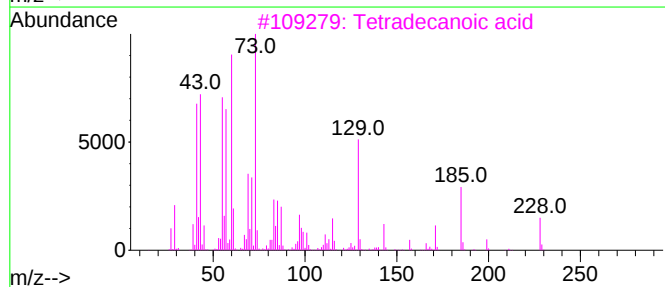
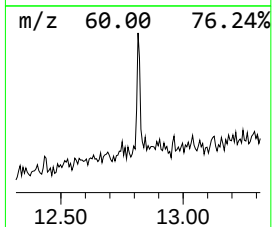
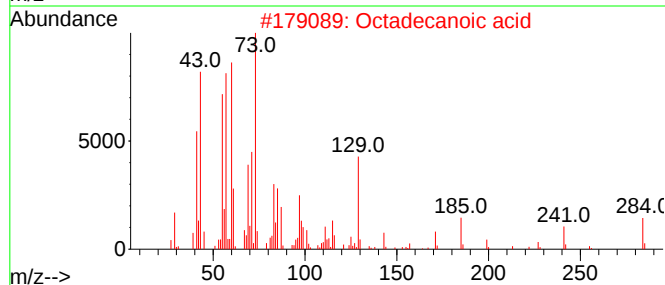
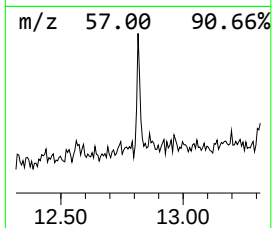
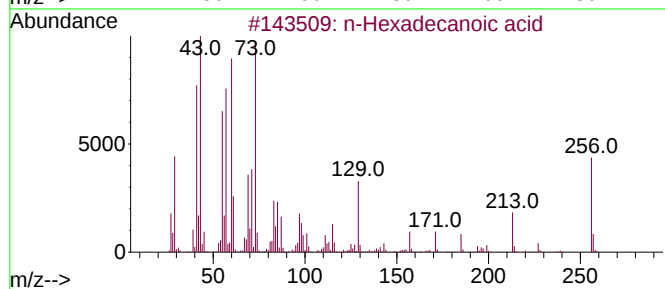
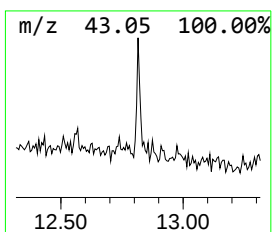
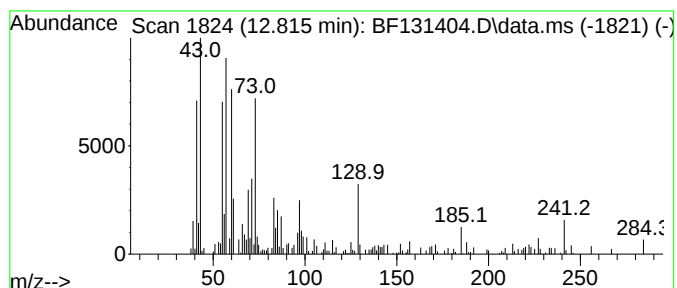
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.815	2.25 ng	78967	Phenanthrene-d10		11.504	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	94
2	Octadecanoic acid		284	C18H36O2	000057-11-4	86
3	Tetradecanoic acid		228	C14H28O2	000544-63-8	86
4	Pentadecanoic acid		242	C15H30O2	001002-84-2	86
5	Tridecanoic acid		214	C13H26O2	000638-53-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-31

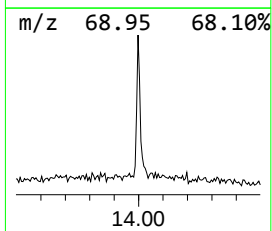
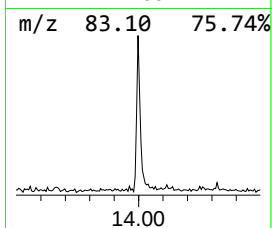
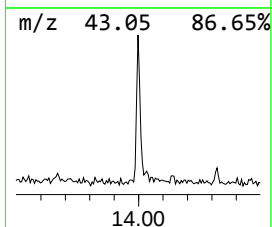
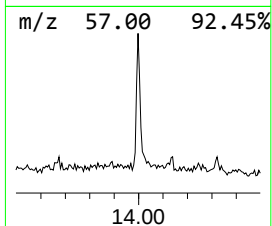
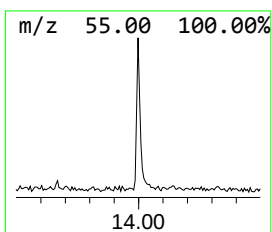
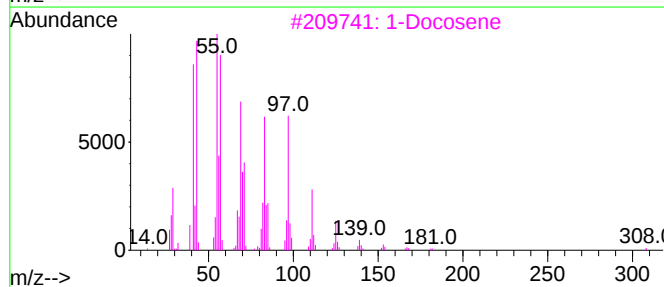
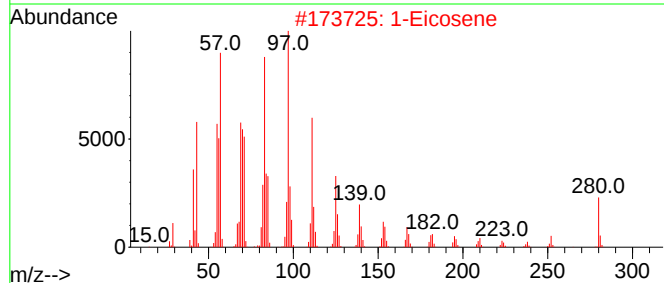
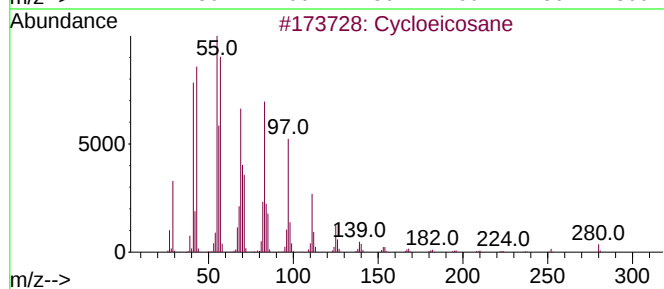
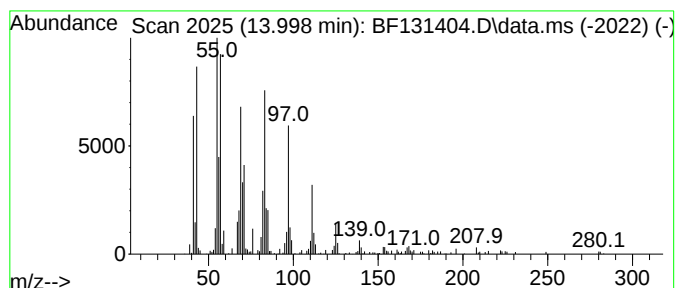
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cycloeicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.998	7.87 ng	189508	Chrysene-d12	14.162

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cycloeicosane	280	C20H40	000296-56-0	99
2		1-Eicosene	280	C20H40	003452-07-1	96
3		1-Docosene	308	C22H44	001599-67-3	94
4		5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-31

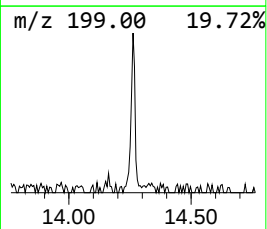
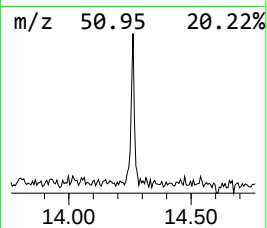
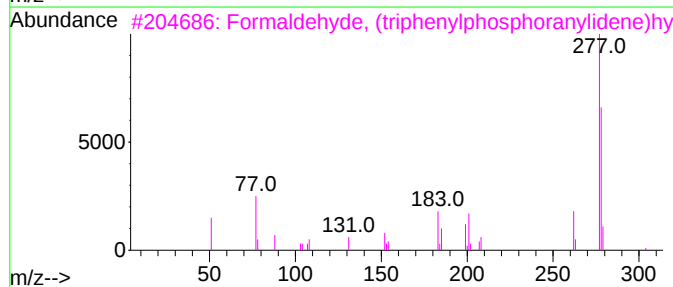
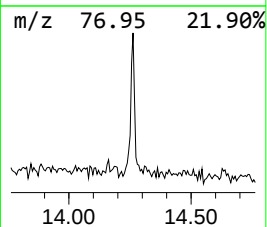
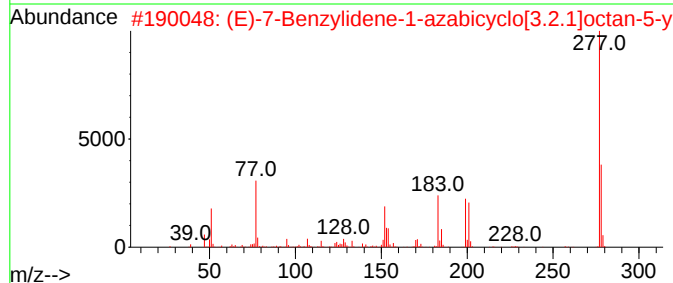
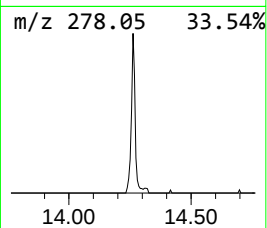
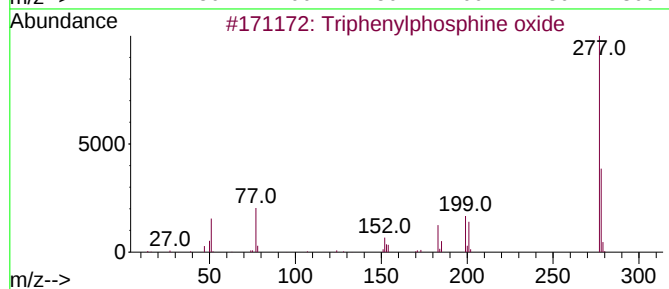
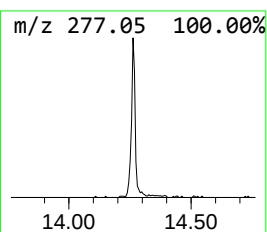
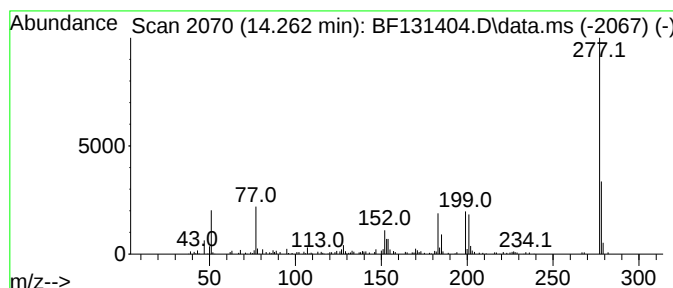
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Triphenylphosphine oxide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.262	3.71 ng	89272	Chrysene-d12	14.162

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	91
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	83
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	45
4			Cobalt,(.eta.<5>-2,4-cyclopentad...	277	C16H15BCo	036682-12-9	37
5			1H-Indole-3-acetic acid, 5-[(tri...	277	C14H19N03Si	072101-35-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
Data File : BF131404.D
Acq On : 25 Nov 2022 19:47
Operator : CG\JU
Sample : N5741-12
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-31

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.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
Butane, 2-metho...	2.251	67.7	ng	1730450	1	6.957	510993	20.0
2-Pentanone, 4-...	5.169	6.7	ng	171241	1	6.957	510993	20.0
Indane	7.133	6.0	ng	152412	1	6.957	510993	20.0
Benzocyclohepta...	9.057	3.7	ng	126440	2	8.245	679861	20.0
Benzoic acid, 2...	9.492	2.4	ng	87370	3	10.010	740474	20.0
unknown9.586	9.586	2.3	ng	85927	3	10.010	740474	20.0
Benzoic acid, p...	9.904	4.8	ng	176831	3	10.010	740474	20.0
Hexathiane	10.321	8.4	ng	312077	3	10.010	740474	20.0
Benzophenone	10.727	2.8	ng	102078	3	10.010	740474	20.0
n-Hexadecanoic ...	12.027	7.9	ng	276819	4	11.504	701776	20.0
Octadecanoic acid	12.815	2.3	ng	78967	4	11.504	701776	20.0
Cycloeicosane	13.998	7.9	ng	189508	5	14.162	481421	20.0
Triphenylphosph...	14.262	3.7	ng	89272	5	14.162	481421	20.0