LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-44

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-BF021622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127175.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.416	356	362	366	rBV2	30020	40660	1.12%	0.179%
2	5.104	475	479	485	rBV	45362	49957	1.37%	0.219%
3	5.510	538	548	555	rBV	1712309	1540864	42.34%	6.768%
4	6.516	715	719	725	rVV	1119459	1003112	27.56%	4.406%
5	6.569	725	728	732	rVB	125374	102306	2.81%	0.449%
6	6.710	749	752	757	rVB	94308	76146	2.09%	0.334%
7	6.886	779	782	785	rBV	786672	628083	17.26%	2.759%
8	6.939	788	791	798	rVB	88742	85813	2.36%	0.377%
9	7.063	810	812	816	rVB	203040	173238	4.76%	0.761%
10	7.128	820	823	833	rVB	145289	144077	3.96%	0.633%
11	7.204	833	836	837	rBV	58788	59256	1.63%	0.260%
12	7.222	837	839	844	rVB	49680	45685	1.26%	0.201%
13	7.275	845	848	857	rBV	92042	103523	2.84%	0.455%
14	7.345	857	860	862	rBV	59425	50184	1.38%	0.220%
15	7.416	868	872	874	rBV	253466	241607	6.64%	1.061%
16	7.457	874	879	882	rVB	2422225	2377555	65.33%	10.443%
17	7.569	894	898	901	rBV2	93345	85791	2.36%	0.377%
18	7.651	909	912	914	rBV	199706	151206	4.15%	0.664%
19	7.681	914	917	919	rVB	167416	148164	4.07%	0.651%
20	7.769	928	932	937	rVB3	45657	53695	1.48%	0.236%
21	7.816	937	940	942	rBV	52554	50512	1.39%	0.222%
22	7.839	942	944	949	rVB2	141108	129810	3.57%	0.570%
23	7.904	949	955	958	rBV	464687	410909	11.29%	1.805%
24	7.981	964	968	970	rBV3	39063	41186	1.13%	0.181%
25	8.010	970	973	975	rVB	43602	39783	1.09%	0.175%
26	8.081	980	985	988	rBV2	50780	85085	2.34%	0.374%
27	8.169	991	1000	1005	rVV	955473	1195581	32.85%	5.252%
28	8.210	1005	1007	1010	rVB	136117	116355	3.20%	0.511%
29	8.286	1017	1020	1024	rBV2	44126	47860	1.32%	0.210%
30	8.351	1027	1031	1036	rVB5	35864	50274	1.38%	0.221%
31	8.445	1044	1047	1052	rVB2	77968	86804	2.39%	0.381%
32	8.528	1056	1061	1063	rBV3	40442	64425	1.77%	0.283%
33	8.551	1063	1065	1068	rVV2	56935	42363	1.16%	0.186%
34	8.633	1076	1079	1083	rBV	43572	42870	1.18%	0.188%
35	8.757	1097	1100	1103	rVB4	40476	46278	1.27%	0.203%
36	8.980	1135	1138	1142	rBV2	41290	49276	1.35%	0.216%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

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-BF021622.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	9.251	1178	1184	1187	rBV	4067418	3639215	100.00%	15.985%
38	9.316	1192	1195	1198	rVB2	54510	51483	1.41%	0.226%
39	9.698	1257	1260	1269	rVB	62615	117649	3.23%	0.517%
40	9.775	1269	1273	1275	rBV4	30088	46636	1.28%	0.205%
41	9.927	1296	1299	1303	rBV	1049607	920301	25.29%	4.042%
42	10.180	1338	1342	1348	rBV	41345	57995	1.59%	0.255%
43	10.463	1388	1390	1394	rBV	52916	49413	1.36%	0.217%
44	10.722	1430	1434	1438	rBV	2530737	2456888	67.51%	10.792%
45	10.827	1450	1452	1461	rVB5	38445	60475	1.66%	0.266%
46	11.057	1488	1491	1497	rBV2	40132	60953	1.67%	0.268%
47	11.363	1540	1543	1546	rVB	112065	88548	2.43%	0.389%
48	11.421	1549	1553	1556	rBV	1143616	908400	24.96%	3.990%
49	11.580	1577	1580	1587	rBV2	75767	98002	2.69%	0.430%
50	11.939	1637	1641	1644	rBV	78208	88440	2.43%	0.388%
51	11.974	1644	1647	1649	rVB	83046	62703	1.72%	0.275%
52	13.010	1818	1823	1826	rBV	3616229	3225076	88.62%	14.166%
53	13.945	1977	1982	1989	rBV4	74036	162319	4.46%	0.713%
54	14.062	1999	2002	2007	rBV	572667	542124	14.90%	2.381%
55	15.562	2251	2257	2264	rVB	363680	469364	12.90%	2.062%

Sum of corrected areas: 22766277

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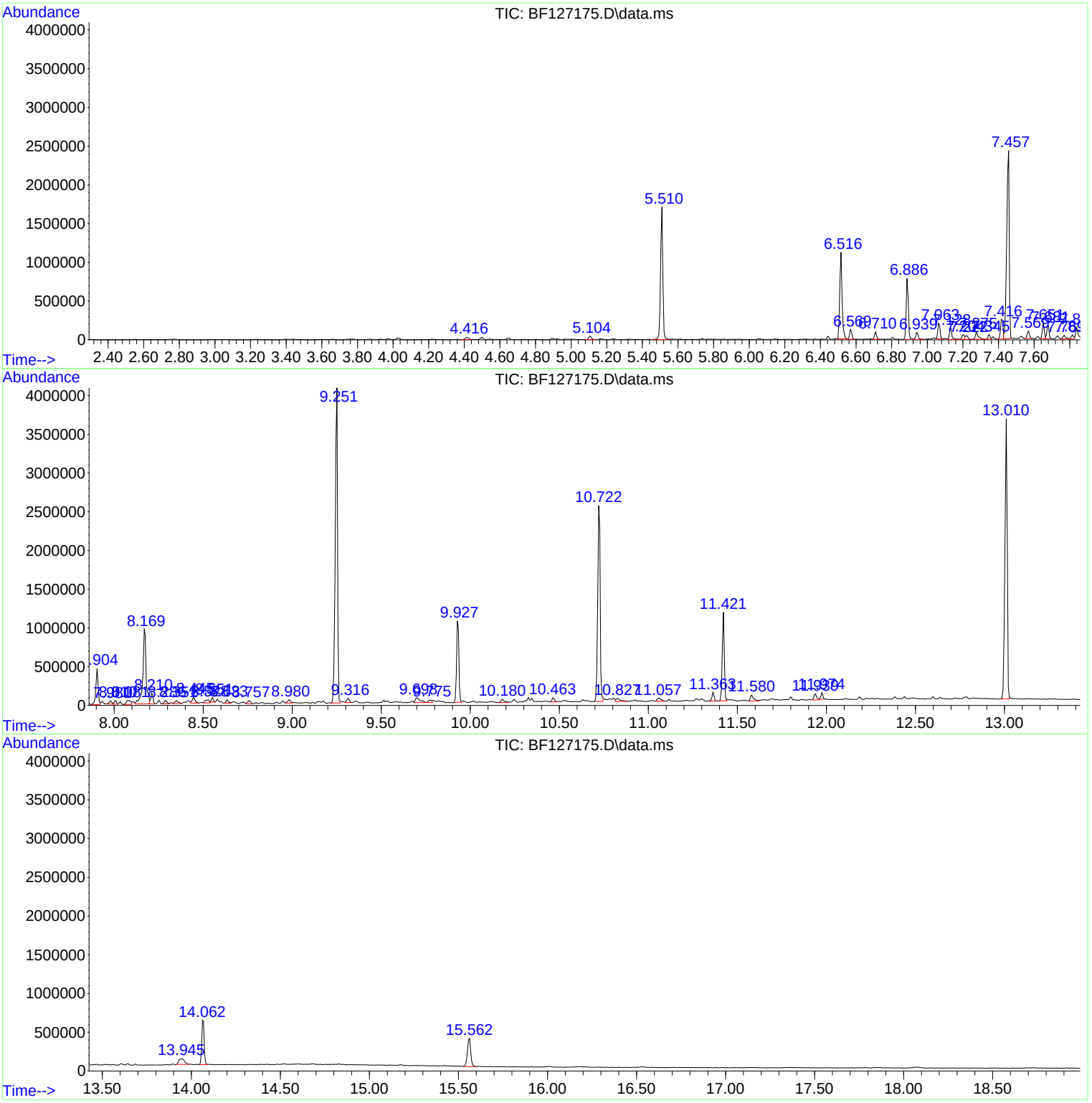
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

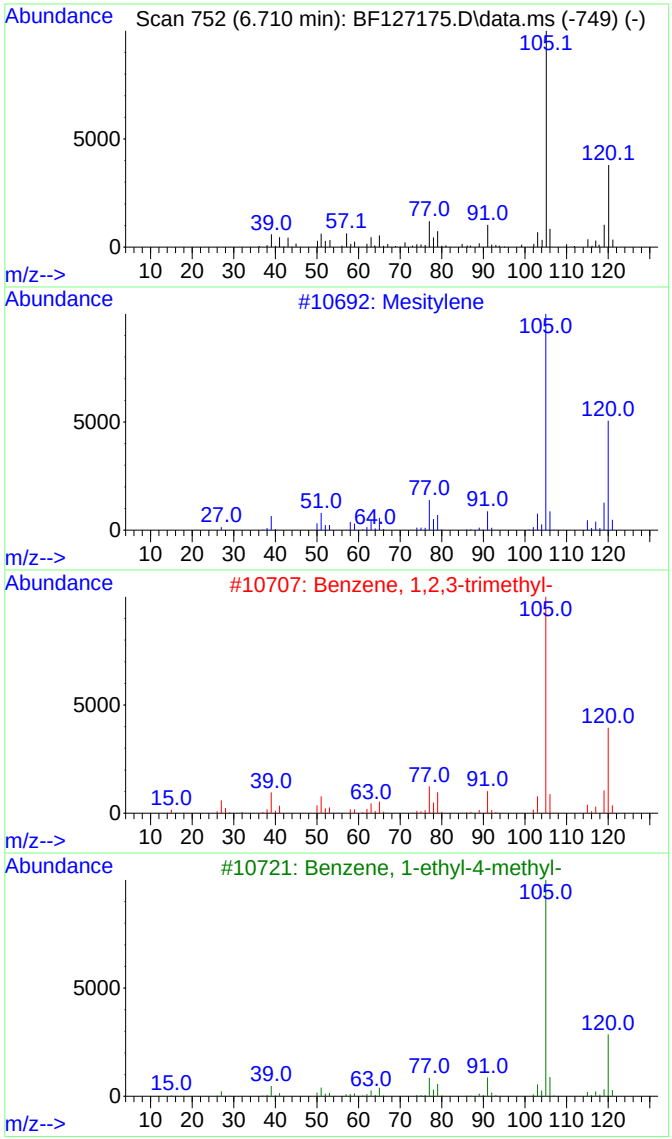
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

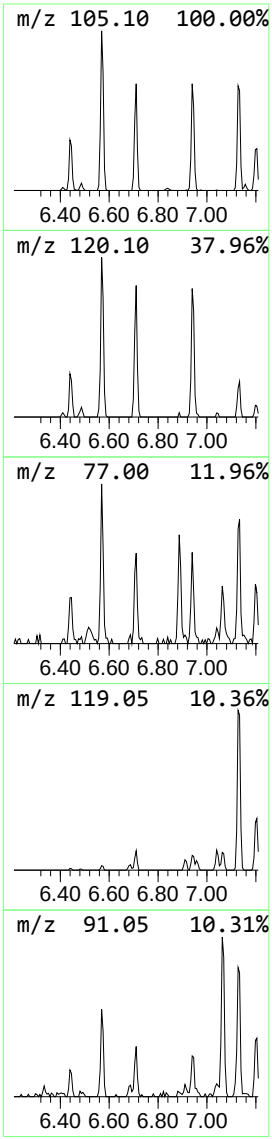
****

Peak Number 1 Mesitylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.710	2.42 ng	76146	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

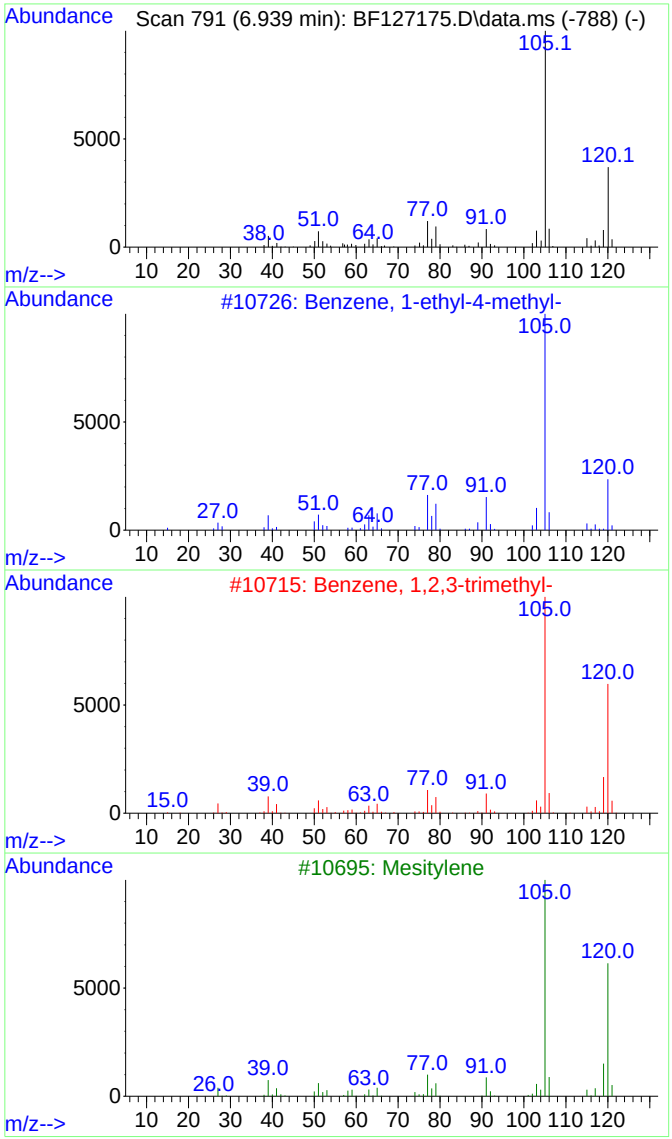
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

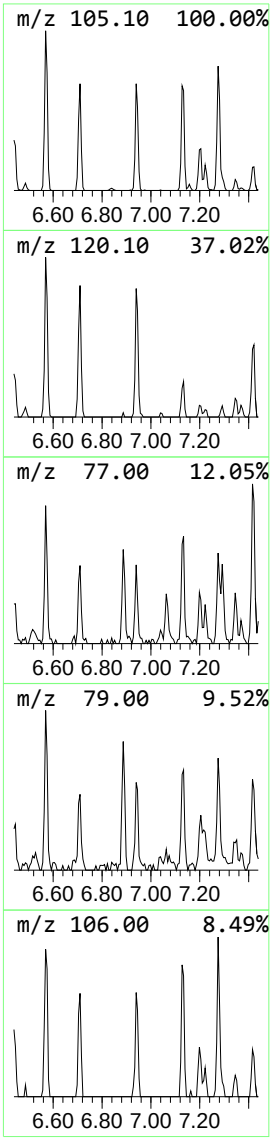
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.939	2.73 ng	85813	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

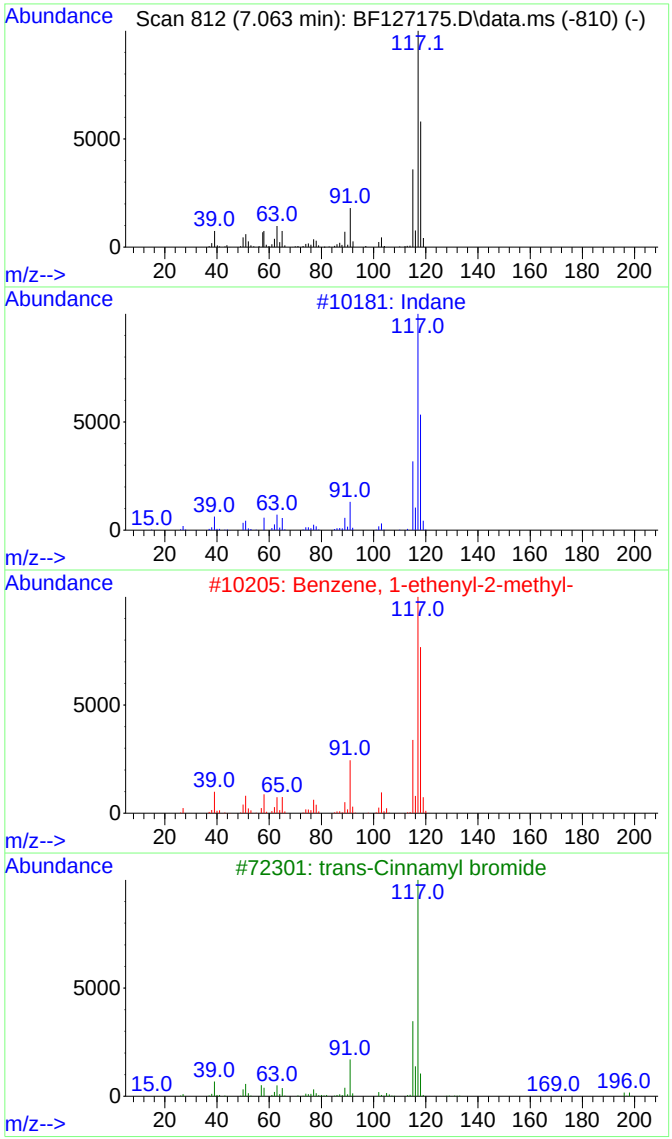
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

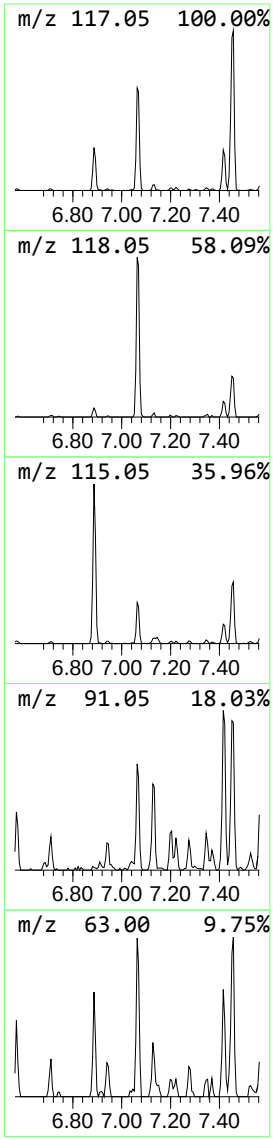
****

Peak Number 3 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.063	5.52 ng	173238	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
3	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	59
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5	Benzene, (isocyanomethyl)-	117	C8H7N	010340-91-7	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

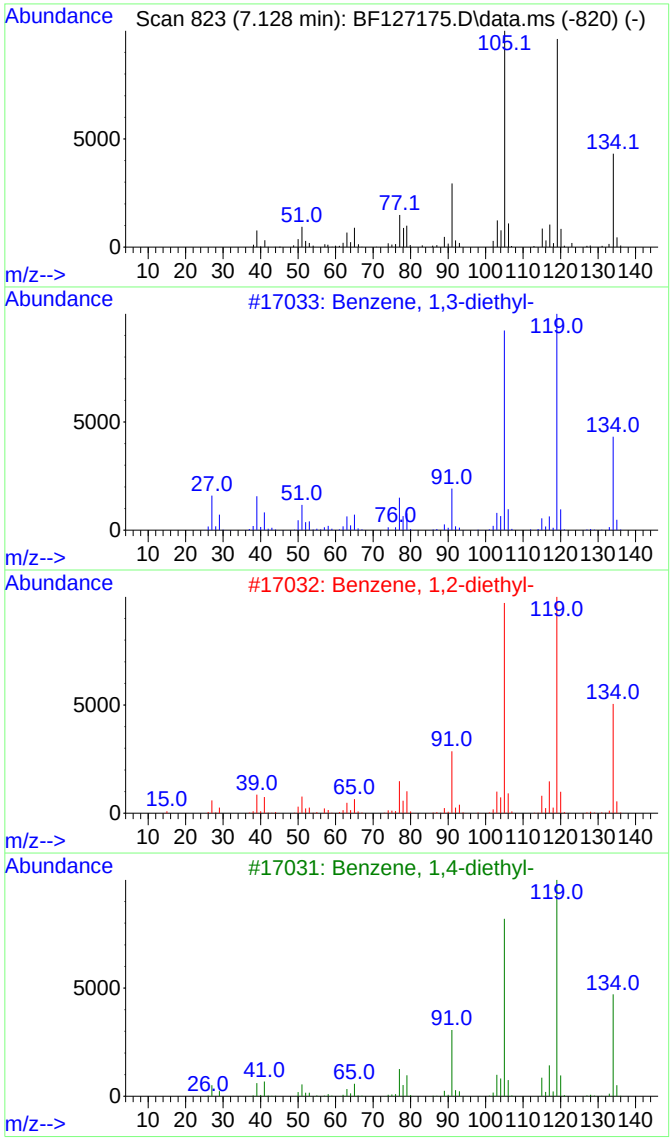
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

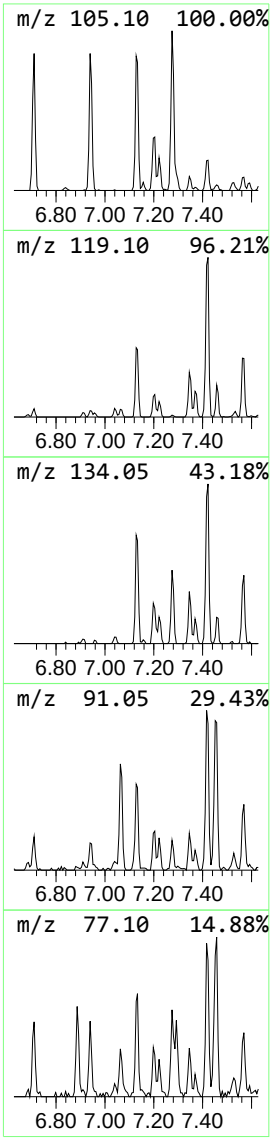
****

Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.128	4.59 ng	144077	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





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Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

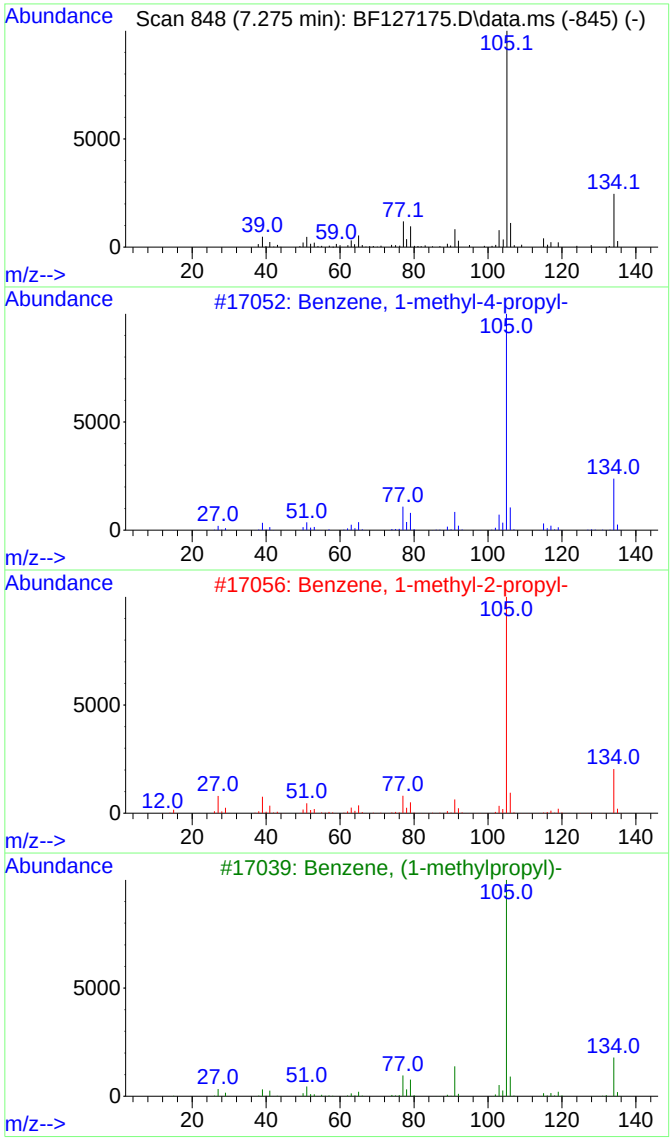
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

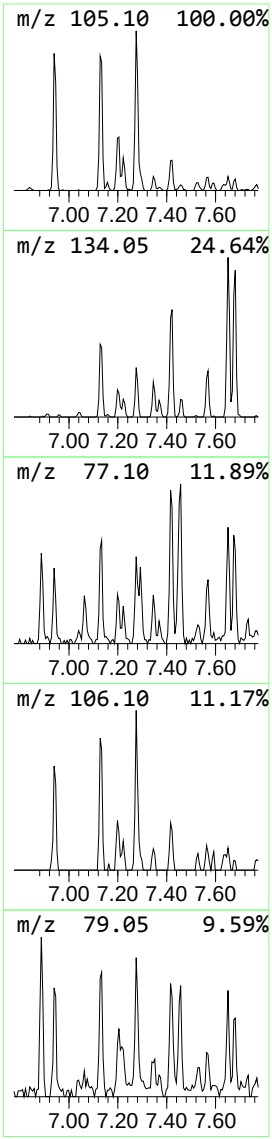
****

Peak Number 5 Benzene, 1-methyl-4-propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.275	3.30 ng	103523	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
5	2-Indanol	134	C9H10O	004254-29-9	90





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Acq On : 03 Mar 2022 17:07
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Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

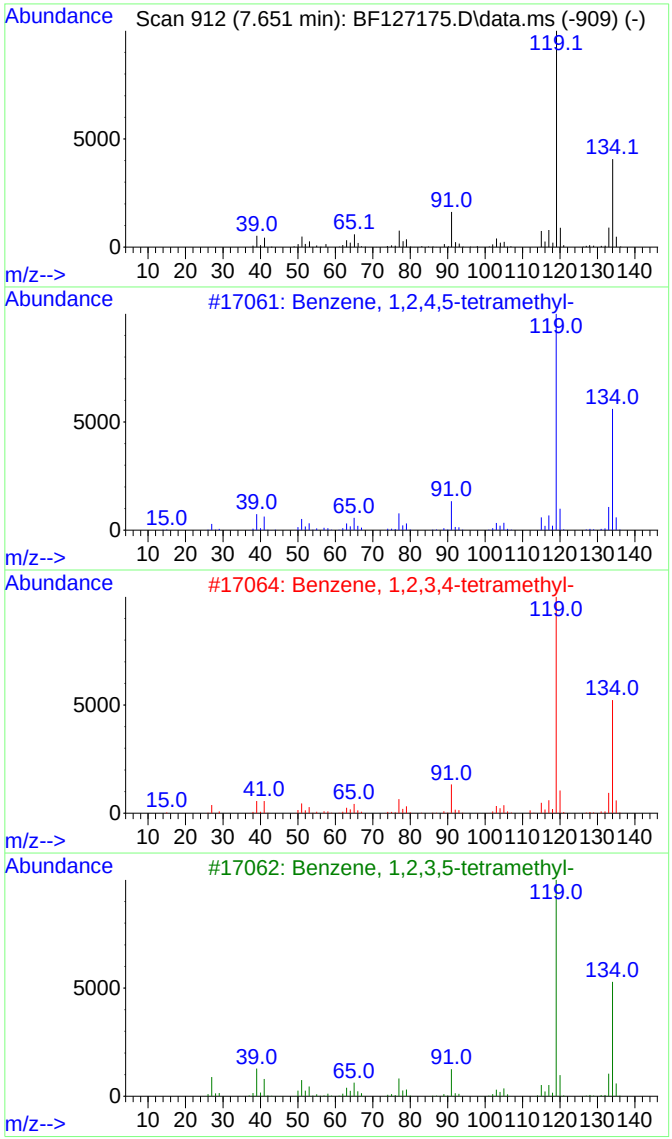
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

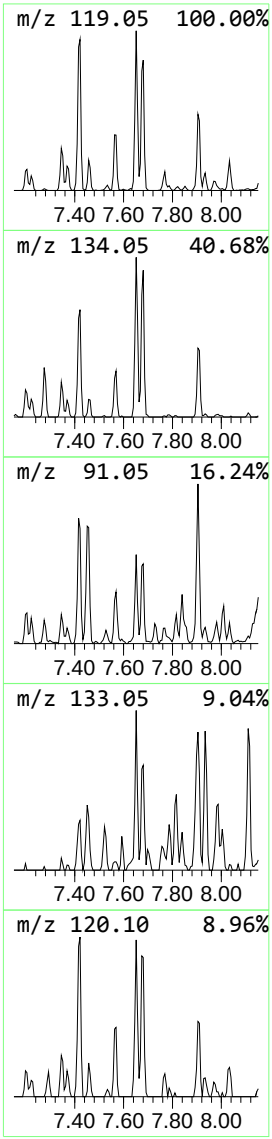
****

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.651	2.53 ng	151206	Naphthalene-d8	8.169

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93





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Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

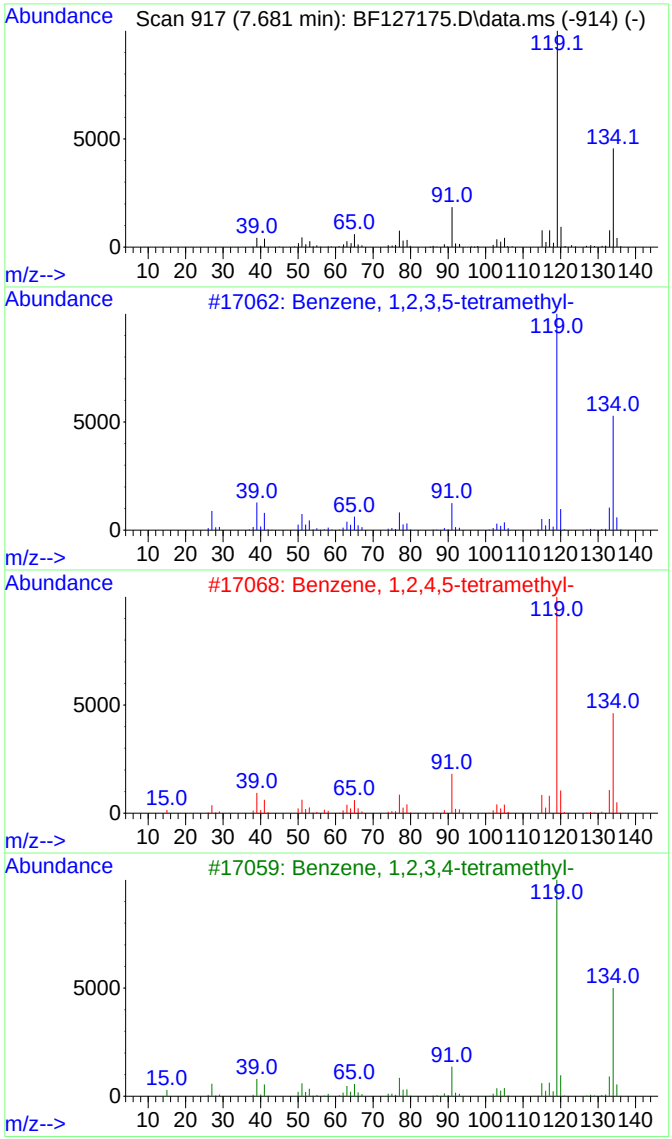
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

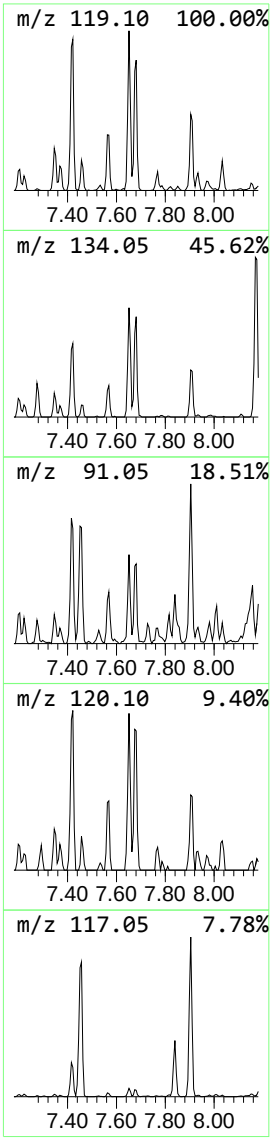
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.681	2.48 ng	148164	Naphthalene-d8	8.169

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

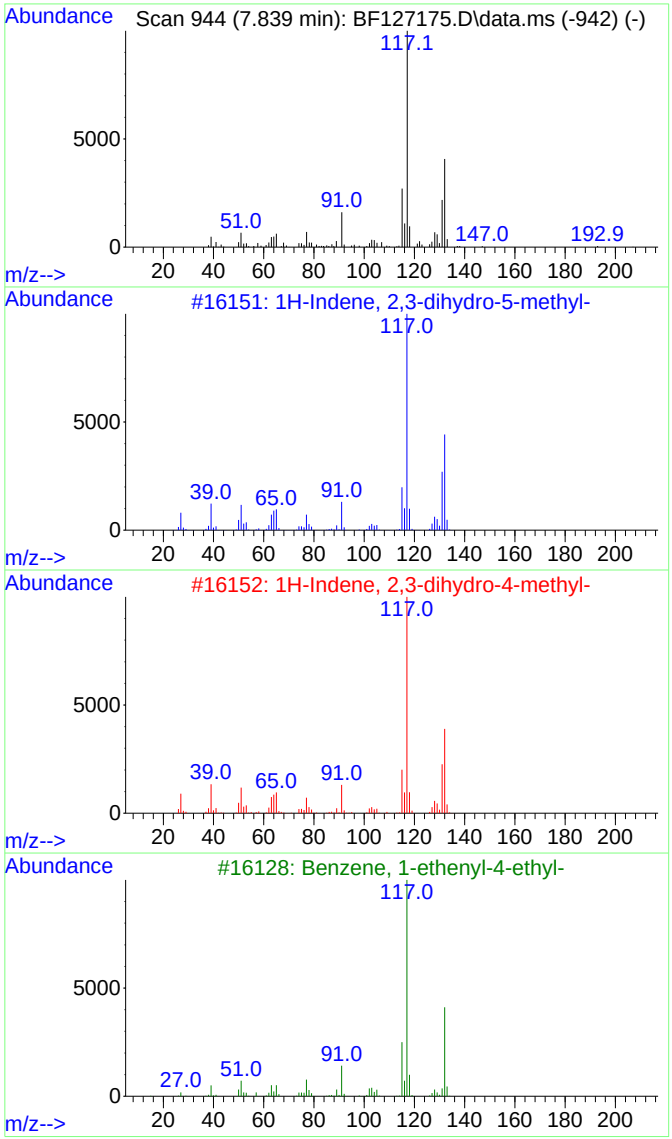
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

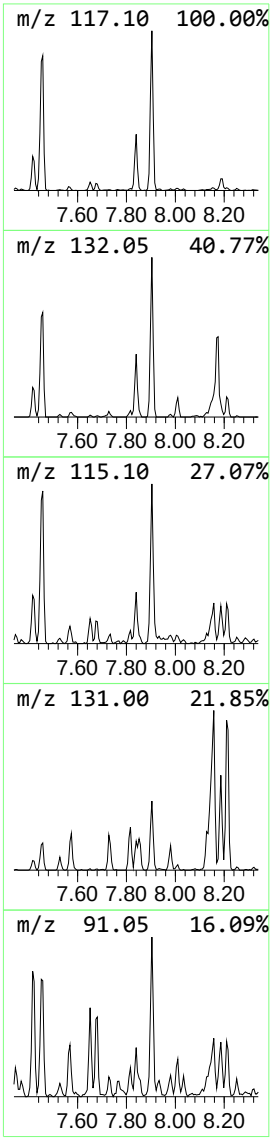
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.839	2.17 ng	129810	Naphthalene-d8	8.169

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

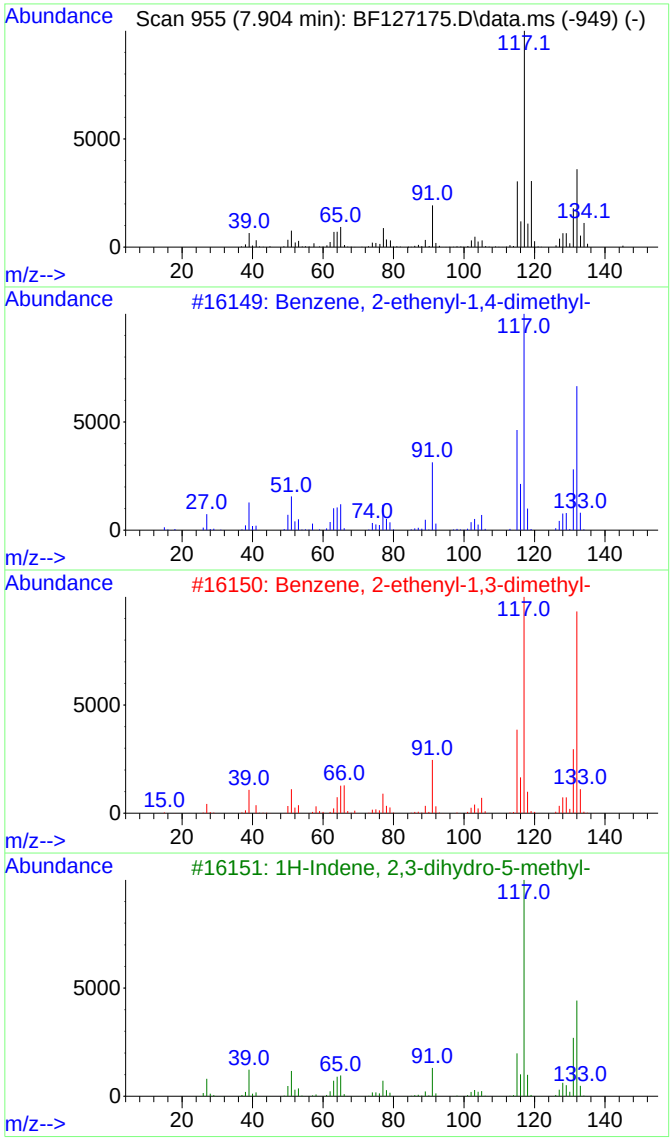
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

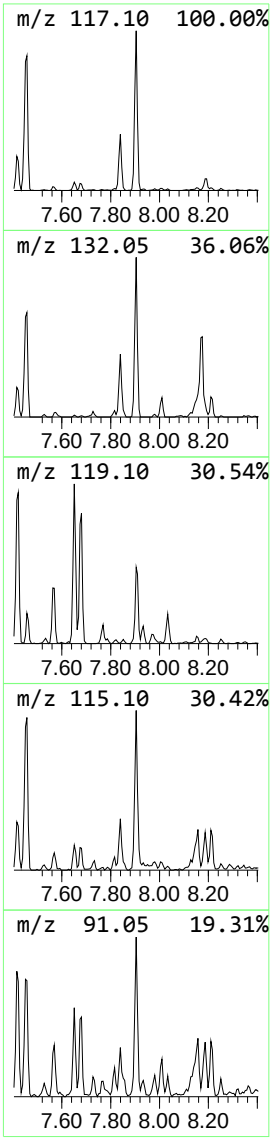
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.904	6.87 ng	410909	Naphthalene-d8	8.169

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





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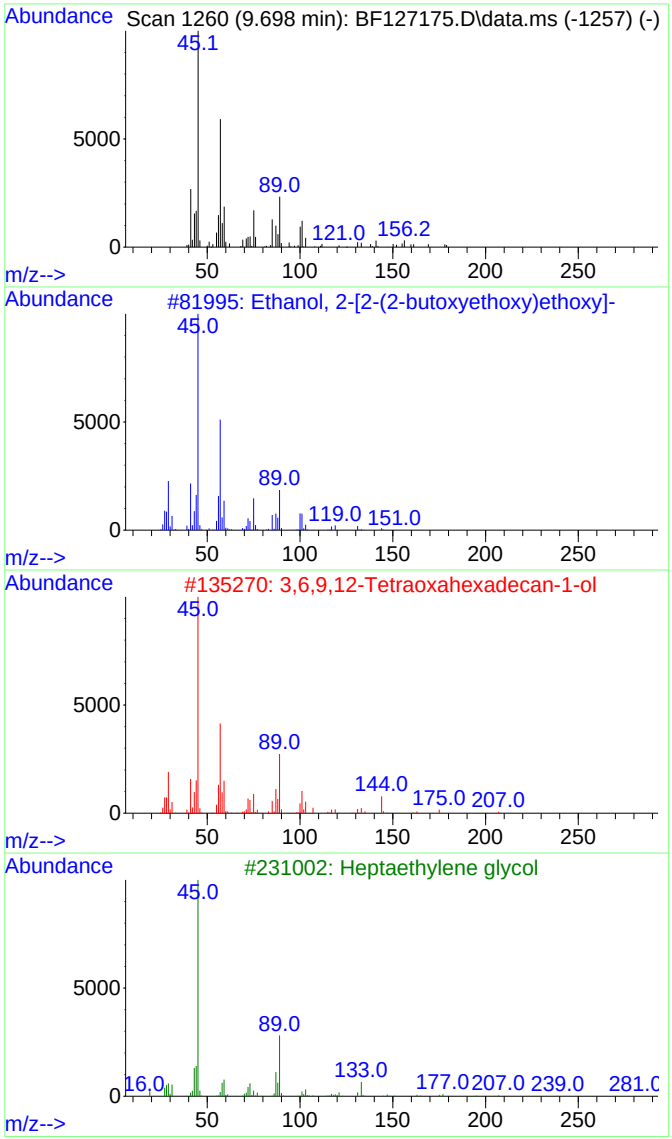
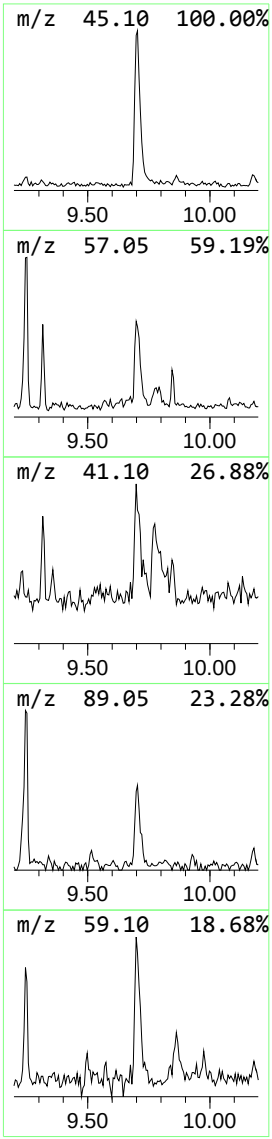
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Ethanol, 2-[2-(2-butoxyetho... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.698	2.56 ng	117649	Acenaphthene-d10	9.927

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	91
2	3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	78
3	Heptaethylene glycol	326	C14H30O8	005617-32-3	64
4	Tetraethylene glycol	194	C8H18O5	000112-60-7	59
5	Hexaethylene glycol	282	C12H26O7	002615-15-8	59





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Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

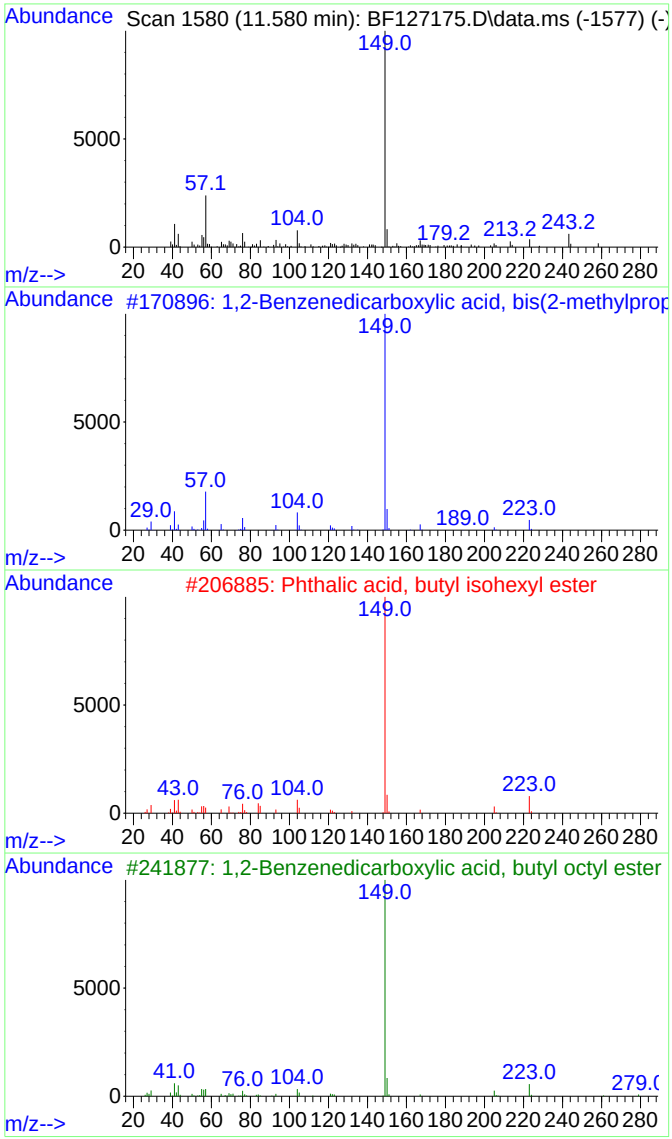
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

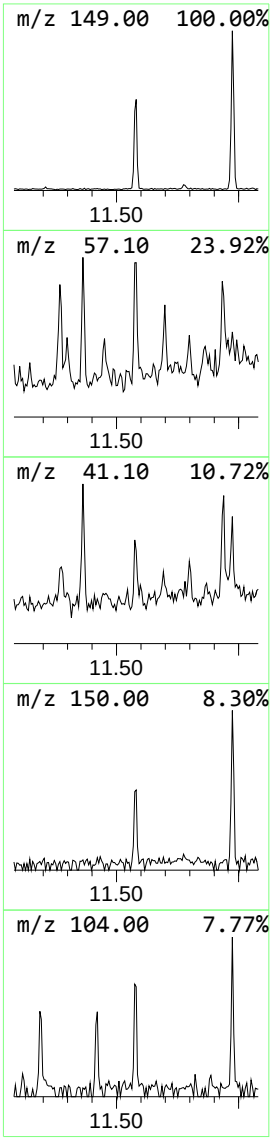
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1,2-Benzenedicarboxylic aci... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.580	2.16 ng	98002	Phenanthrene-d10	11.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	80
2	Phthalic acid, butyl isohexyl ester	306	C18H26O4	1000309-03-6	72
3	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000084-78-6	72
4	Phthalic acid, hept-4-yl isobuty...	320	C19H28O4	1000356-78-3	72
5	Phthalic acid, 2-fluorophenyl he...	344	C20H21FO4	1000360-51-6	72





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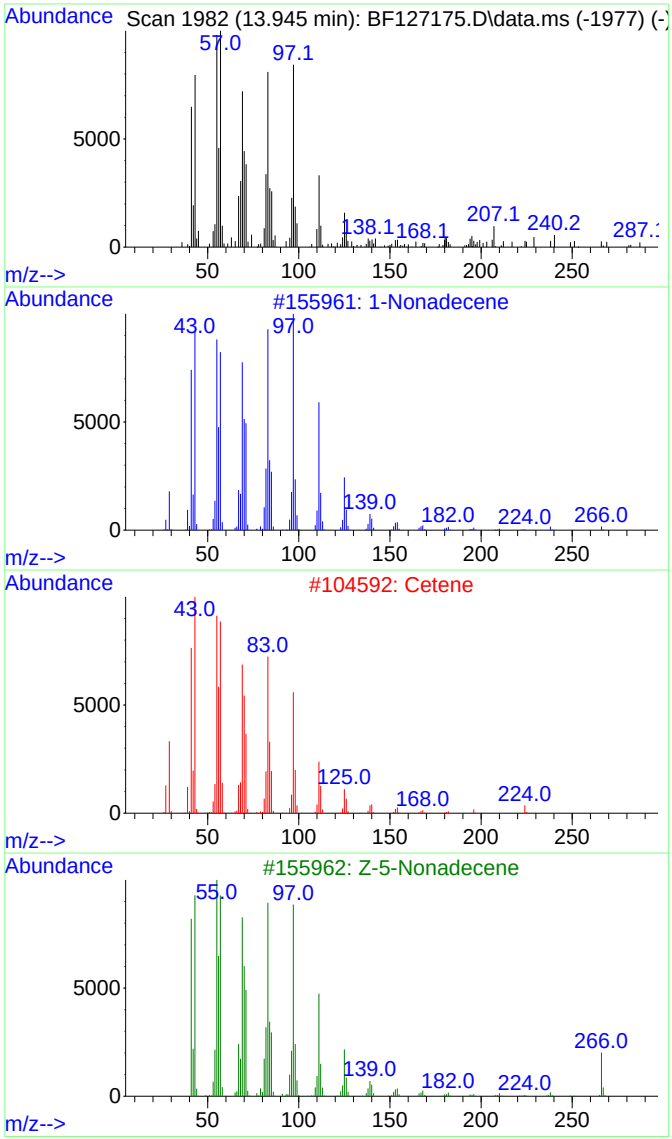
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

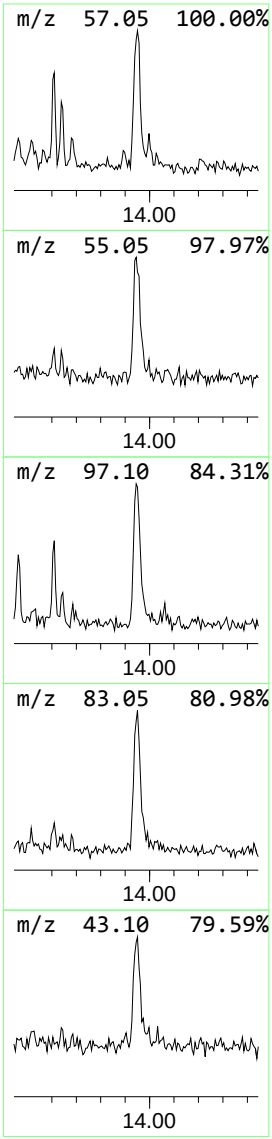
****

Peak Number 12 1-Nonadecene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.945	5.99 ng	162319	Chrysene-d12	14.068

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	99
2	Cetene	224	C16H32	000629-73-2	98
3	Z-5-Nonadecene	266	C19H38	1000131-11-8	95
4	1-Heptadecene	238	C17H34	006765-39-5	95
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127175.D
Acq On : 03 Mar 2022 17:07
Operator : CG\JU
Sample : N1689-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-44

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Mesitylene	6.710	2.4	ng	76146	1	6.886	628083	20.0
Benzene, 1-ethy...	6.939	2.7	ng	85813	1	6.886	628083	20.0
Indane	7.063	5.5	ng	173238	1	6.886	628083	20.0
Benzene, 1,3-di...	7.128	4.6	ng	144077	1	6.886	628083	20.0
Benzene, 1-meth...	7.275	3.3	ng	103523	1	6.886	628083	20.0
Benzene, 1,2,4,...	7.651	2.5	ng	151206	2	8.169	1195580	20.0
Benzene, 1,2,3,...	7.681	2.5	ng	148164	2	8.169	1195580	20.0
1H-Indene, 2,3-...	7.839	2.2	ng	129810	2	8.169	1195580	20.0
Benzene, 2-ethe...	7.904	6.9	ng	410909	2	8.169	1195580	20.0
Ethanol, 2-[2-(...	9.698	2.6	ng	117649	3	9.927	920301	20.0
1,2-Benzenedica...	11.580	2.2	ng	98002	4	11.422	908400	20.0
1-Nonadecene	13.945	6.0	ng	162319	5	14.068	542124	20.0