LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

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-BP040921.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005288.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.281	367	373	379	rVB	1241672	1815520	55.82%	7.660%
2	6.828	631	636	638	rBV	106611	168265	5.17%	0.710%
3	6.875	638	644	645	rVV	1163062	1915331	58.89%	8.081%
4	6.899	645	648	663	rVB	1597204	2450997	75.36%	10.342%
5	7.299	712	716	725	rVB3	18242	36577	1.12%	0.154%
6	7.675	771	780	787	rBV	240782	380166	11.69%	1.604%
7	7.740	787	791	798	rVV	130199	182634	5.62%	0.771%
8	8.840	970	978	985	rBV	788967	1244168	38.26%	5.250%
9	9.404	1063	1074	1082	rVB5	11097	39370	1.21%	0.166%
10	10.299	1210	1226	1241	rBV	363402	1284382	39.49%	5.419%
11	10.463	1243	1254	1260	rVV	289446	486590	14.96%	2.053%
12	11.487	1423	1428	1437	rBV5	26079	58203	1.79%	0.246%
13	11.987	1506	1513	1524	rBV	566121	897120	27.58%	3.785%
14	12.863	1656	1662	1666	rVB	33893	51392	1.58%	0.217%
15	12.951	1667	1677	1686	rVV	1598552	2190198	67.34%	9.241%
16	13.745	1806	1812	1816	rVV3	38586	56027	1.72%	0.236%
17	13.816	1816	1824	1828	rVV4	59641	126031	3.88%	0.532%
18	13.863	1828	1832	1840	rVB10	17441	36436	1.12%	0.154%
19	14.157	1878	1882	1887	rBV4	34203	49414	1.52%	0.208%
20	14.234	1892	1895	1901	rVB3	31390	45934	1.41%	0.194%
21	14.339	1902	1913	1920	rVB2	441759	696114	21.40%	2.937%
22	15.604	2124	2128	2132	rBV2	58059	80553	2.48%	0.340%
23	15.839	2163	2168	2174	rVB	325602	451484	13.88%	1.905%
24	16.086	2203	2210	2215	rBV2	146306	271430	8.35%	1.145%
25	16.186	2223	2227	2231	rBV	51370	79809	2.45%	0.337%
26	16.245	2233	2237	2243	rVV5	44793	83628	2.57%	0.353%
27	16.304	2245	2247	2251	rVV3	28596	35529	1.09%	0.150%
28	16.875	2339	2344	2348	rBV3	245802	385761	11.86%	1.628%
29	17.104	2378	2383	2387	rBV	579665	844913	25.98%	3.565%
30	17.198	2395	2399	2402	rBV	139025	170216	5.23%	0.718%
31	18.151	2558	2561	2568	rVB	295967	366918	11.28%	1.548%
32	18.992	2700	2704	2713	rVB	279342	411784	12.66%	1.737%
33	19.139	2725	2729	2733	rVB	175025	234669	7.22%	0.990%
34	19.480	2784	2787	2795	rVB	165466	266524	8.19%	1.125%
35	19.680	2816	2821	2828	rVB	2734815	3252296	100.00%	13.723%
36	20.916	3028	3031	3039	rVB3	329541	497626	15.30%	2.100%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

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-BP040921.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	20.986	3040	3043	3050	rVB	196511	254678	7.83%	1.075%
38	21.116	3062	3065	3070	rVB	245157	271138	8.34%	1.144%
39	21.204	3075	3080	3084	rBV	712635	943393	29.01%	3.981%
40	23.474	3461	3466	3473	rVB2	317695	587127	18.05%	2.477%

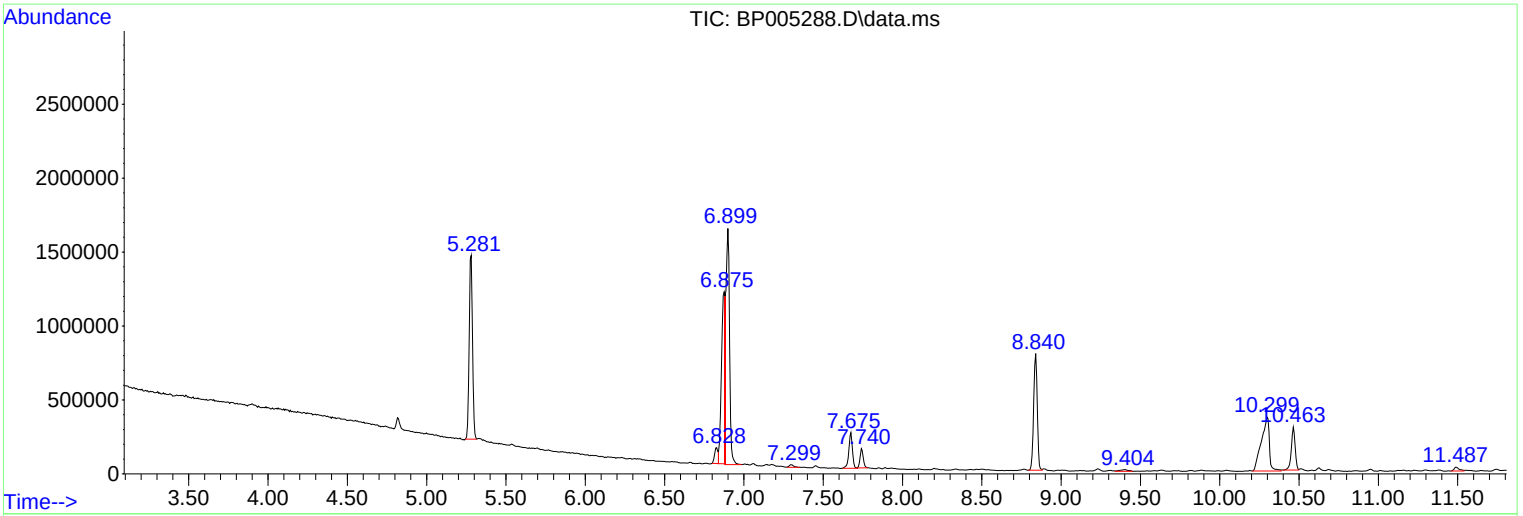
Sum of corrected areas: 23700345

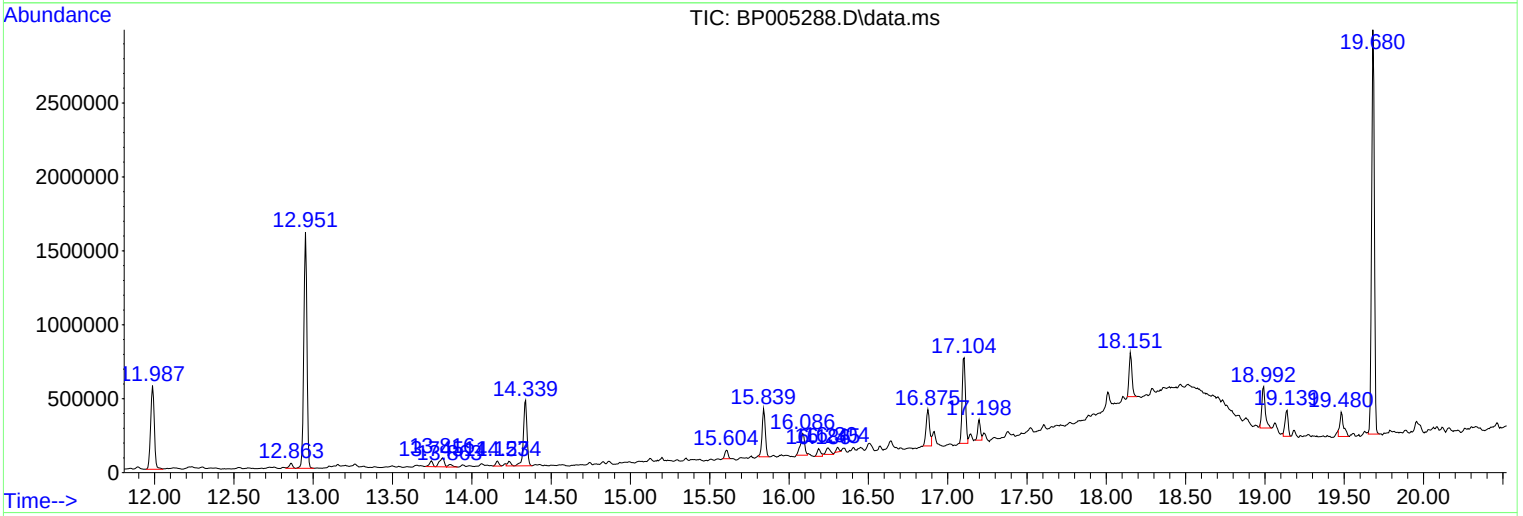
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

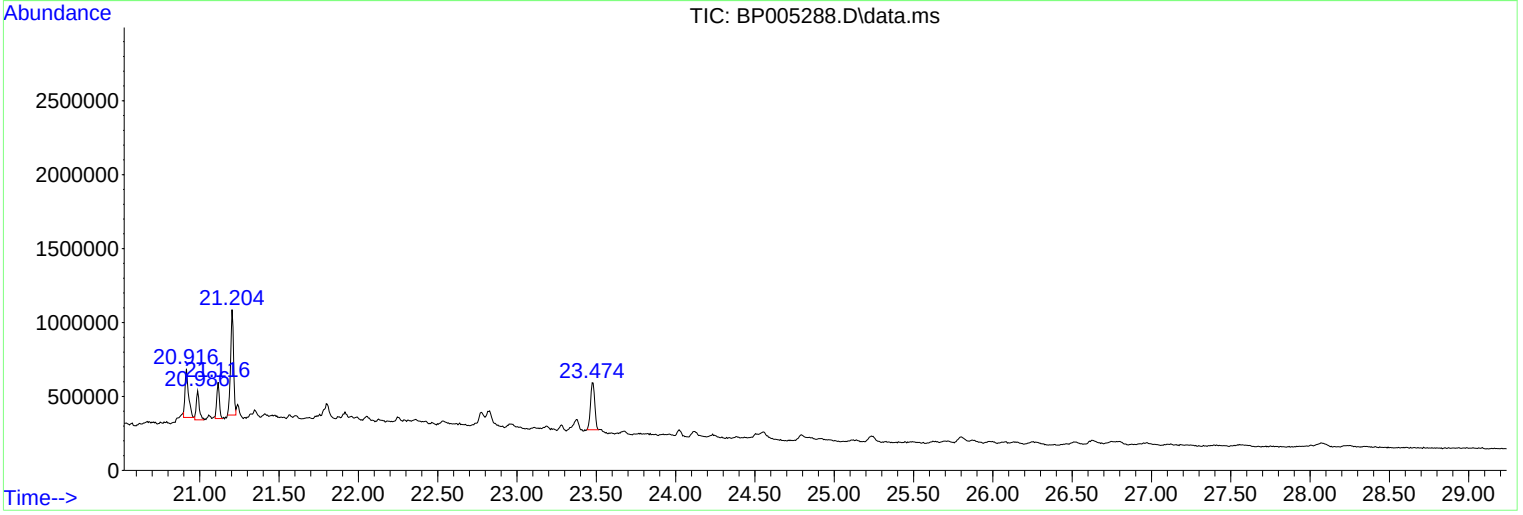
Instrument :
BNA_P
ClientSampleId :
SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

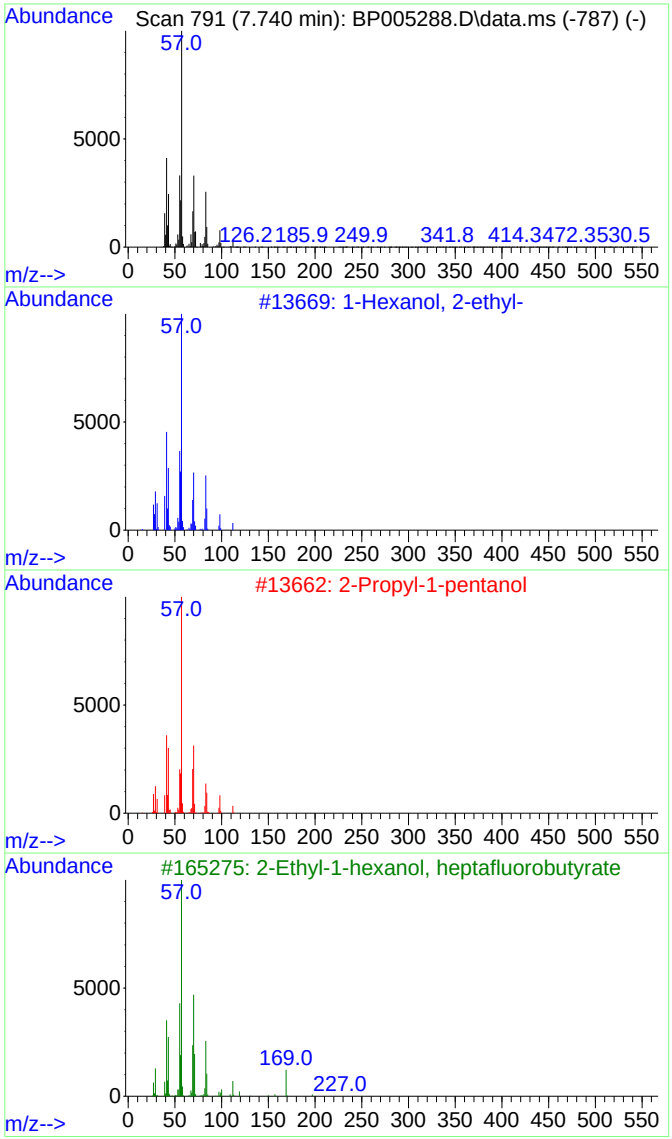
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

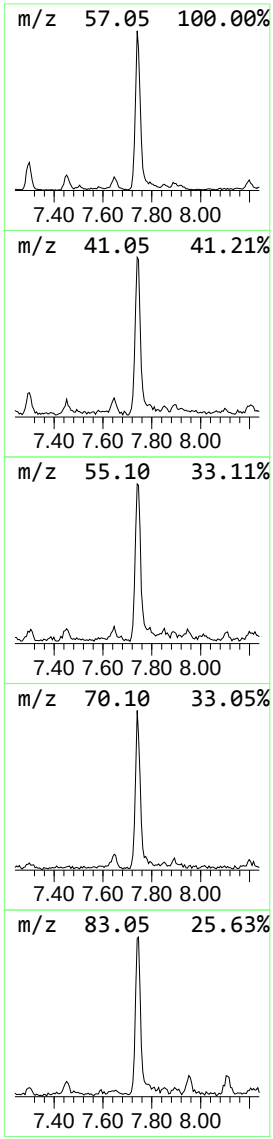
****

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.740	9.61 ng	182634	1,4-Dichlorobenzene-d4	7.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	2-Propyl-1-pentanol	130	C8H18O	058175-57-8	72
3	2-Ethyl-1-hexanol, heptafluorobu...	326	C12H17F7O2	1000365-16-0	64
4	2-Ethyl-1-hexanol, pentafluoropr...	276	C11H17F5O2	1000365-51-1	59
5	2-Ethylhexyl hydrogen maleate	228	C12H20O4	002370-71-0	56





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.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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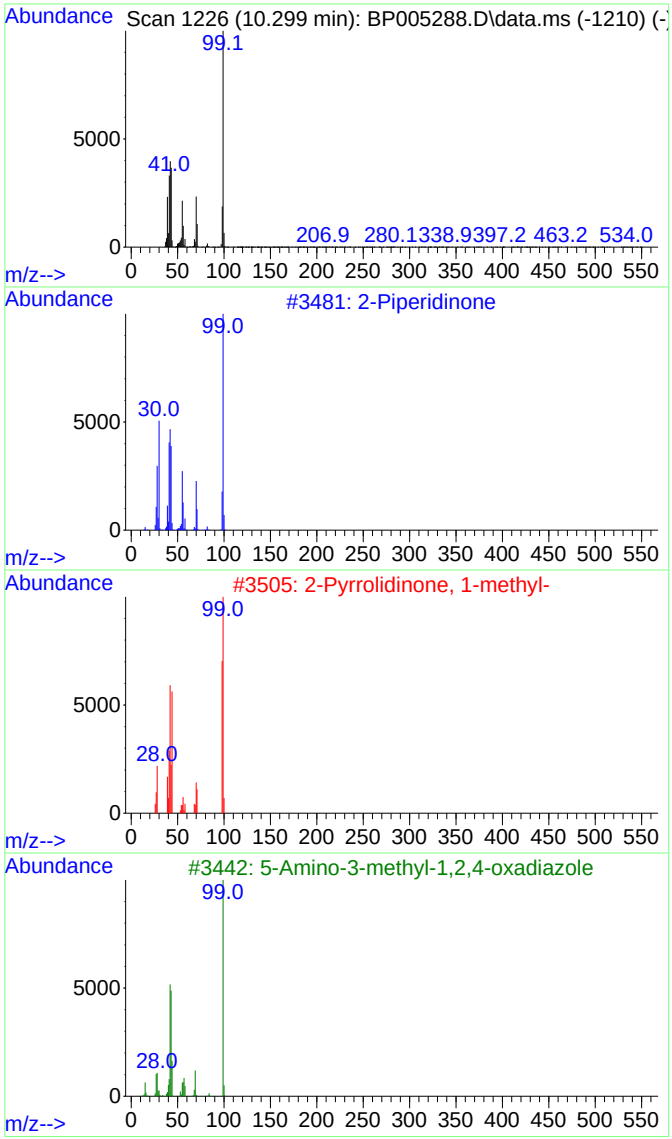
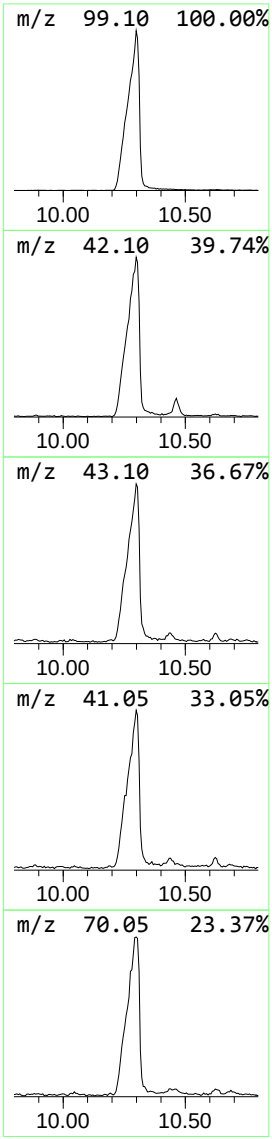
Peak Number 2 2-Piperidinone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.299	52.79 ng	1284380	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Piperidinone	99	C5H9NO	000675-20-7	91
2	2-Pyrrolidinone, 1-methyl-	99	C5H9NO	000872-50-4	72
3	5-Amino-3-methyl-1,2,4-oxadiazole	99	C3H5N3O	003663-39-6	59
4	3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	43
5	Allyl Isothiocyanate	99	C4H5NS	000057-06-7	32





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

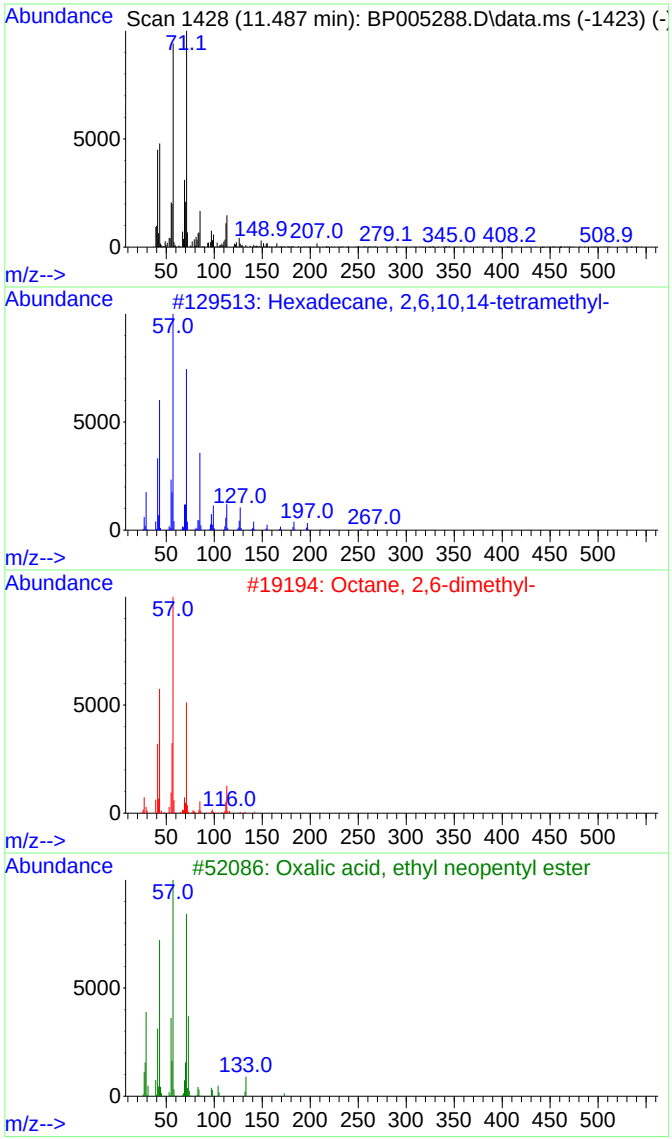
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

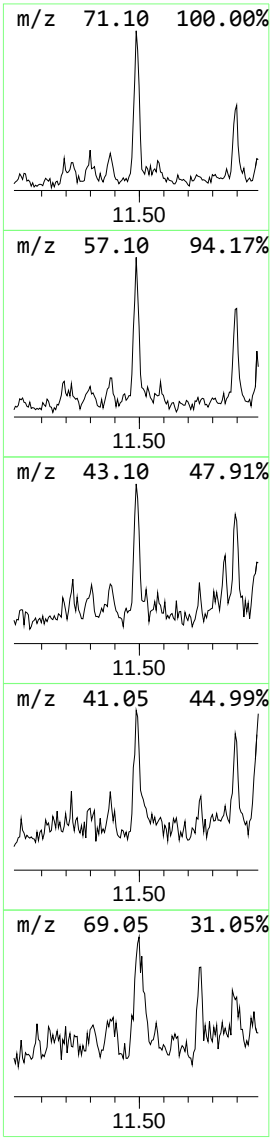
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Hexadecane, 2,6,10,14-tetra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.487	2.39 ng	58203	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
3	Oxalic acid, ethyl neopentyl ester	188	C9H16O4	1000309-72-4	59
4	Sulfurous acid, dodecyl 2-pentyl...	320	C17H36O3S	1000309-16-1	53
5	Nonane, 2,3-dimethyl-	156	C11H24	002884-06-2	53





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.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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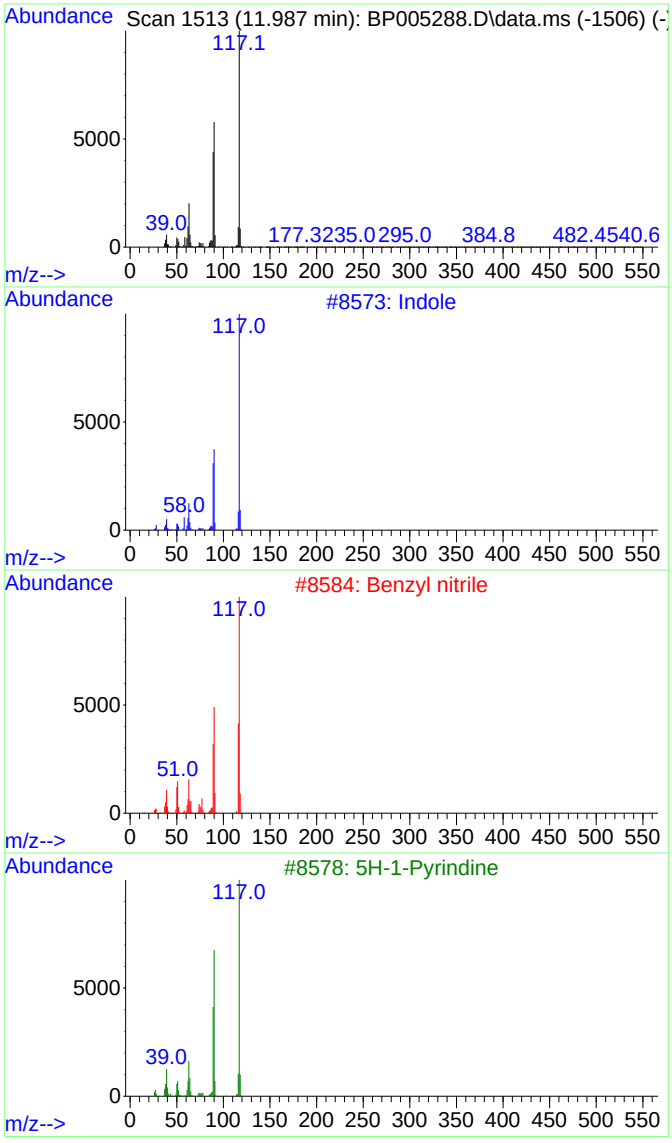
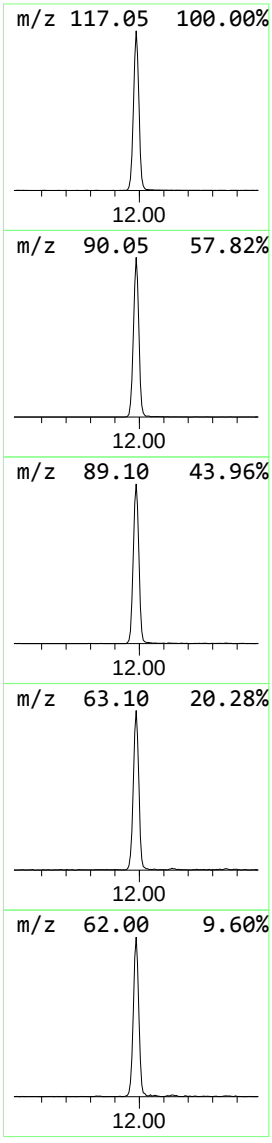
Peak Number 4 Indole Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.987	36.87 ng	897120	Naphthalene-d8	10.463

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole	117	C8H7N	000120-72-9	95
2	Benzyl nitrile	117	C8H7N	000140-29-4	91
3	5H-1-Pyrindine	117	C8H7N	000270-91-7	91
4	Indolizine	117	C8H7N	000274-40-8	91
5	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

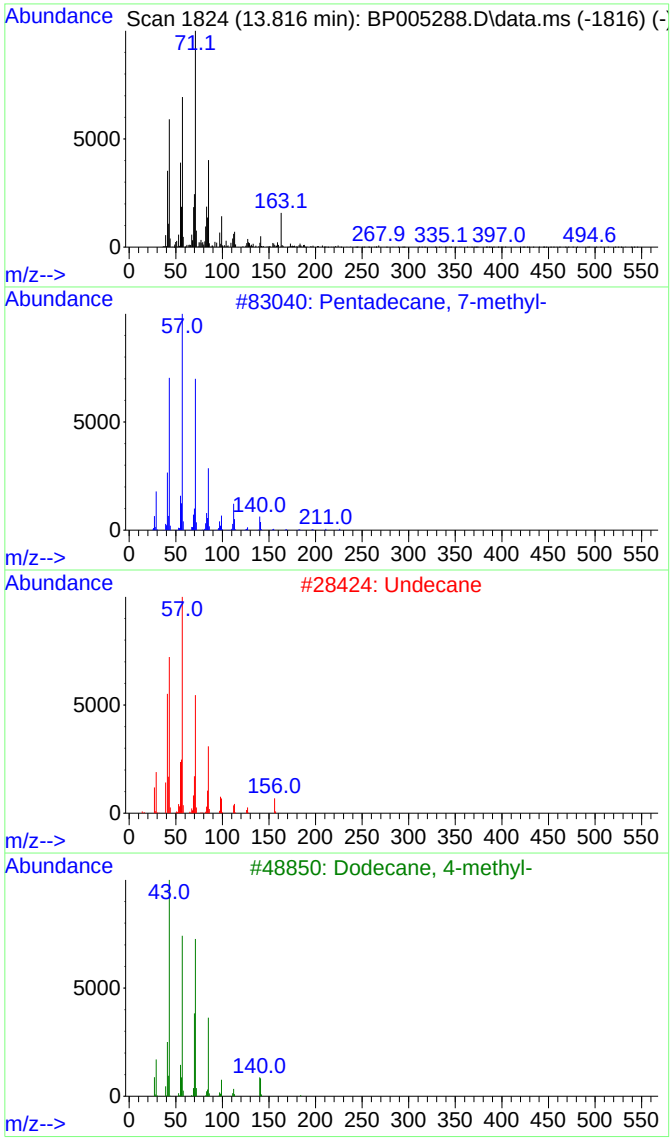
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

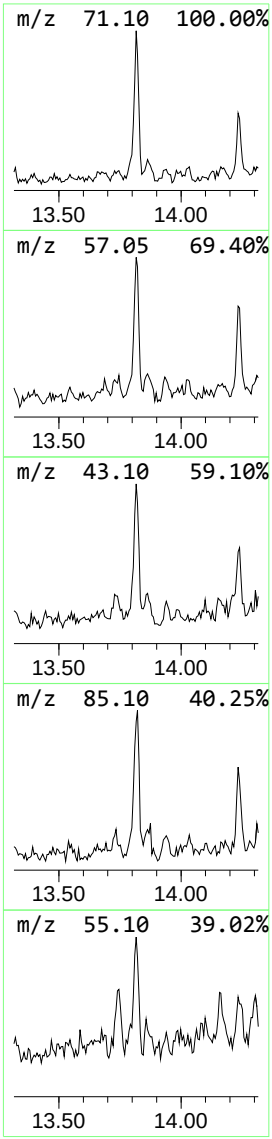
****

Peak Number 5 Pentadecane, 7-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	3.62 ng	126031	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72
2	Undecane	156	C11H24	001120-21-4	58
3	Dodecane, 4-methyl-	184	C13H28	006117-97-1	58
4	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	53
5	Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

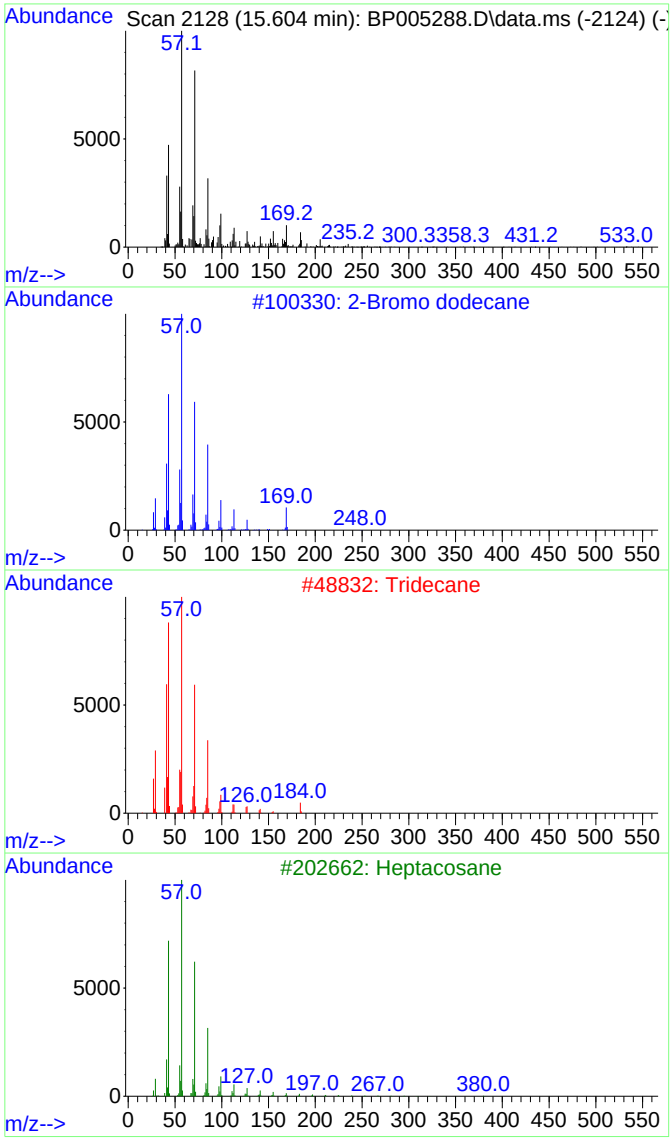
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

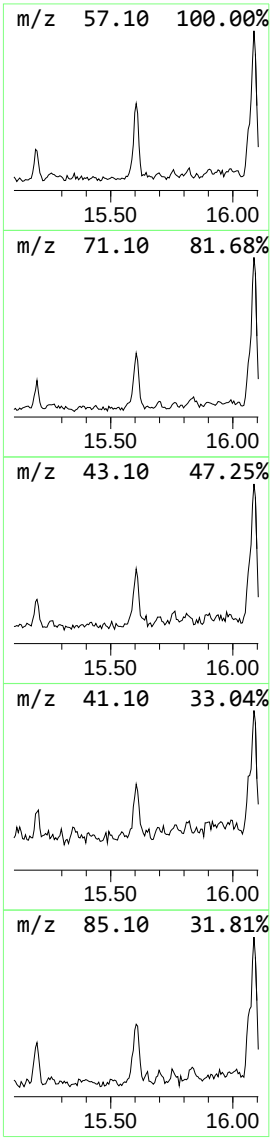
****

Peak Number 6 2-Bromo dodecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.604	2.31 ng	80553	Acenaphthene-d10	14.339

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	89
2	Tridecane	184	C13H28	000629-50-5	80
3	Heptacosane	380	C27H56	000593-49-7	72
4	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	64
5	Tetracosane	338	C24H50	000646-31-1	59





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Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.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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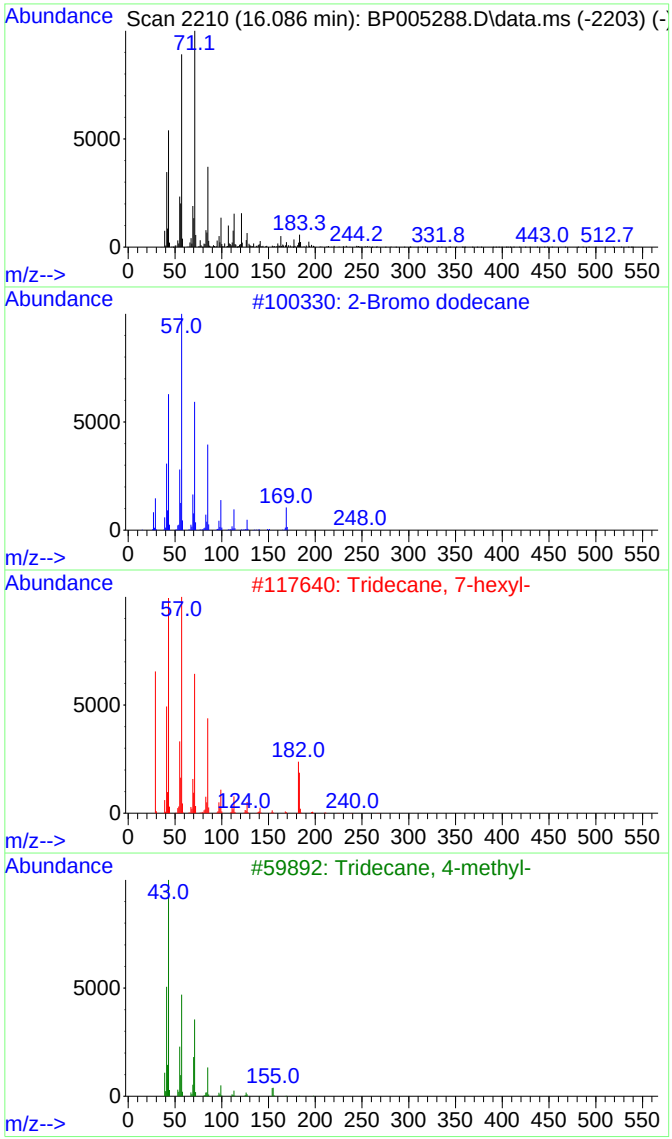
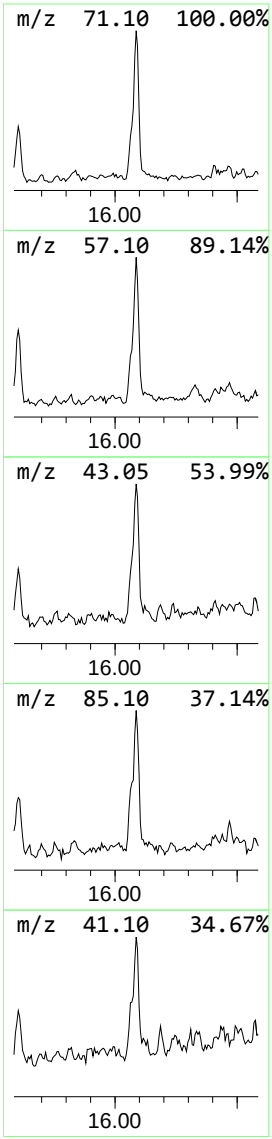
Peak Number 7 Tridecane, 7-hexyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.086	6.43 ng	271430	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane	248	C12H25Br	013187-99-0	92
2	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	78
3	Tridecane, 4-methyl-	198	C14H30	026730-12-1	72
4	Tridecane, 5-propyl-	226	C16H34	055045-11-9	70
5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.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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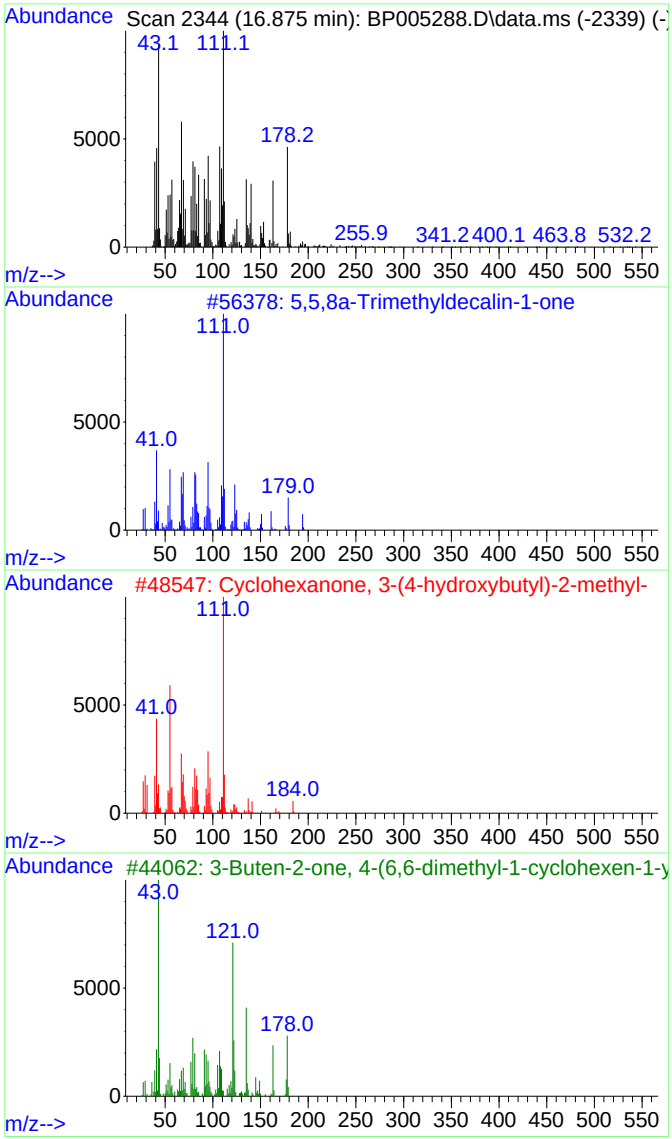
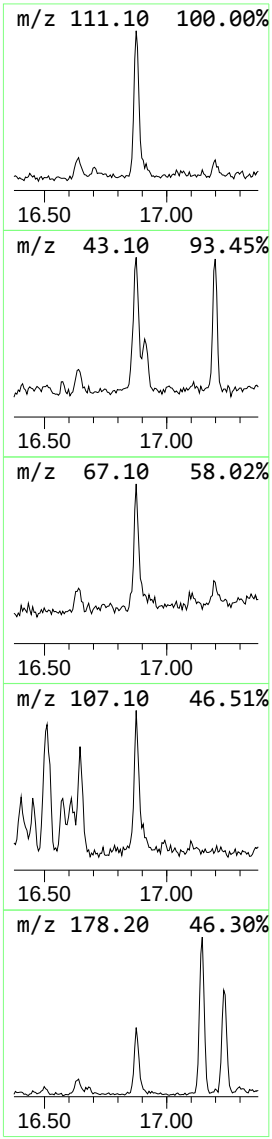
Peak Number 8 unknown16.875 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.875	9.13 ng	385761	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5,8a-Trimethyldecalin-1-one	194	C13H22O	1000195-96-1	43
2	Cyclohexanone, 3-(4-hydroxybutyl...	184	C11H20O2	091212-98-5	35
3	3-Buten-2-one, 4-(6,6-dimethyl-1...	178	C12H18O	065133-79-1	14
4	2-(Cyanoimino)oxazolidine	111	C4H5N3O	050411-06-8	14
5	Tetrahydrocarvone	154	C10H18O	000499-70-7	14





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Data File : BP005288.D
Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

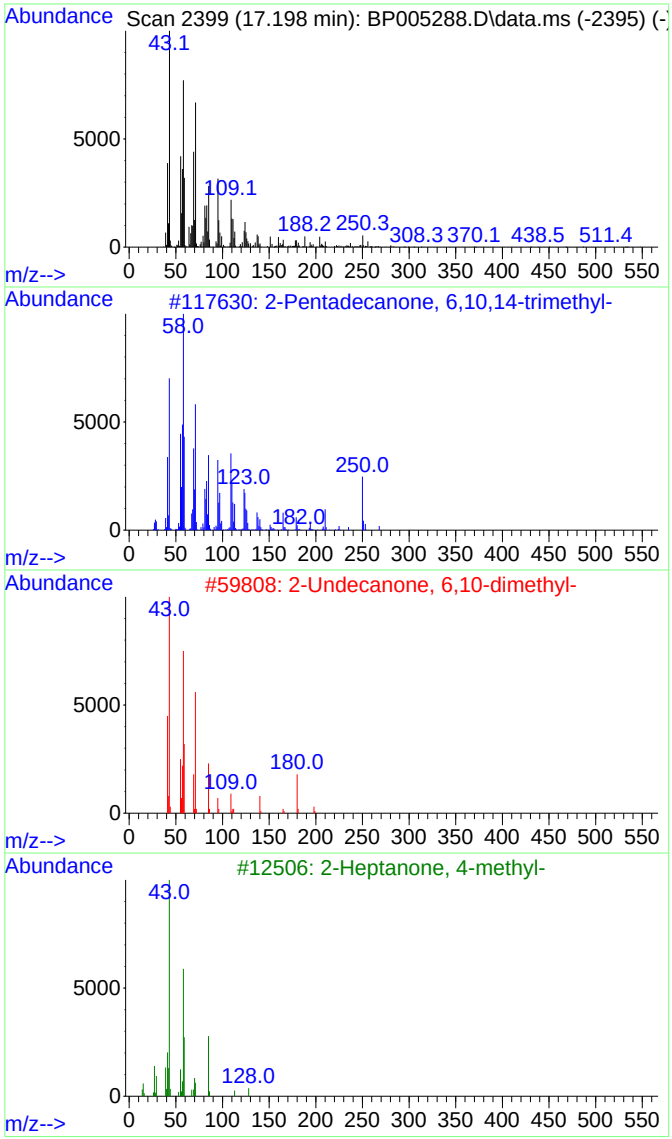
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

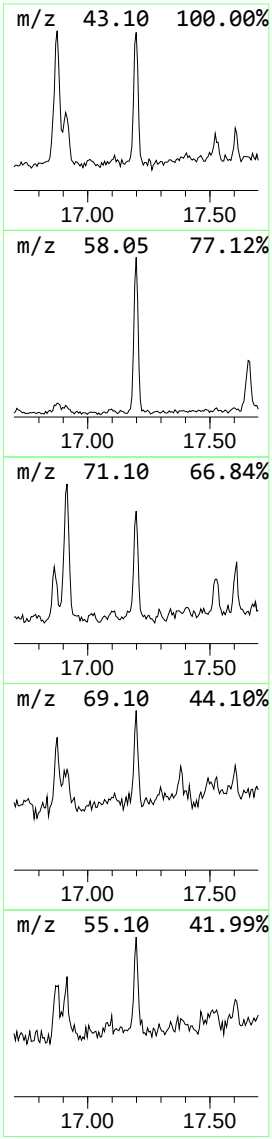
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 2-Pentadecanone, 6,10,14-tr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.198	4.03 ng	170216	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentadecanone, 6,10,14-trimethyl-	268	C18H36O	000502-69-2	93
2	2-Undecanone, 6,10-dimethyl-	198	C13H26O	001604-34-8	59
3	2-Heptanone, 4-methyl-	128	C8H16O	006137-06-0	43
4	Cyclopentolate	291	C17H25NO3	000512-15-2	35
5	1,3-Propanediamine, N,N,N',N'-te...	130	C7H18N2	000110-95-2	35





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Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

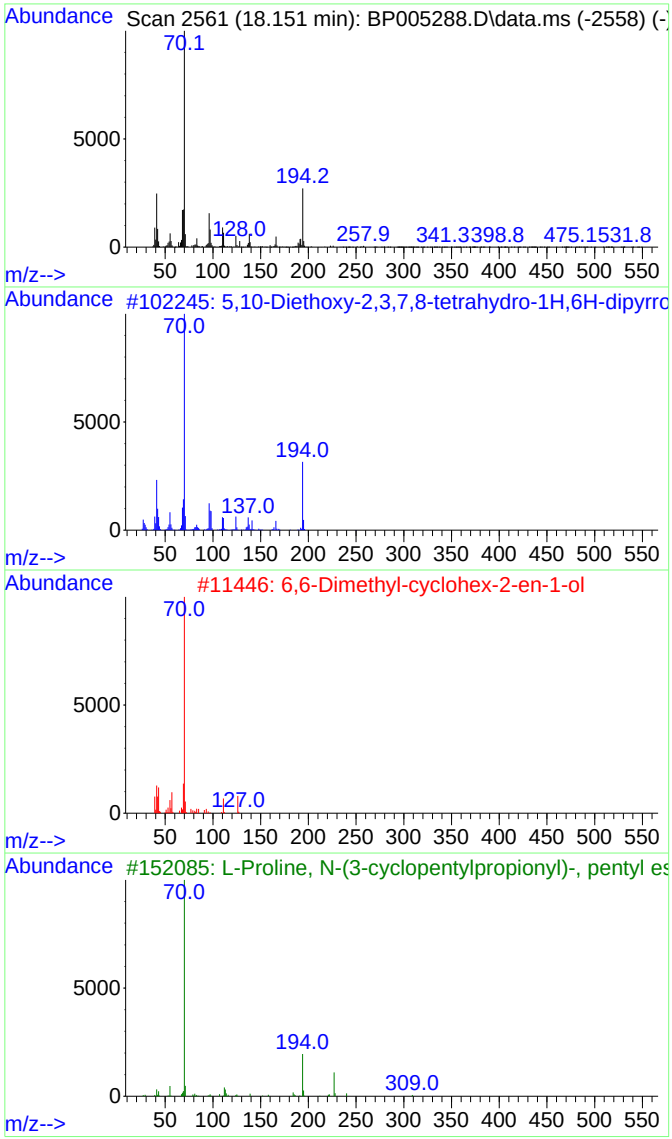
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

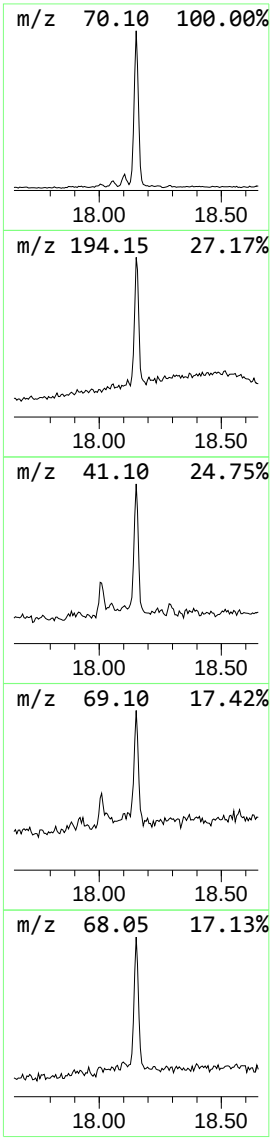
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 5,10-Diethoxy-2,3,7,8-tetra... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	8.69 ng	366918	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,10-Diethoxy-2,3,7,8-tetrahydro...	250	C14H22N2O2	1000190-75-5	64
2	6,6-Dimethyl-cyclohex-2-en-1-ol	126	C8H14O	038313-10-9	43
3	L-Proline, N-(3-cyclopentylpropi...	309	C18H31NO3	1000346-41-0	40
4	2,7-Dimethyl-2,7-octadien-1-amine	153	C10H19N	077984-58-8	36
5	L-Proline, N-(3-cyclopentylpropi...	295	C17H29NO3	1000346-40-9	33





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Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
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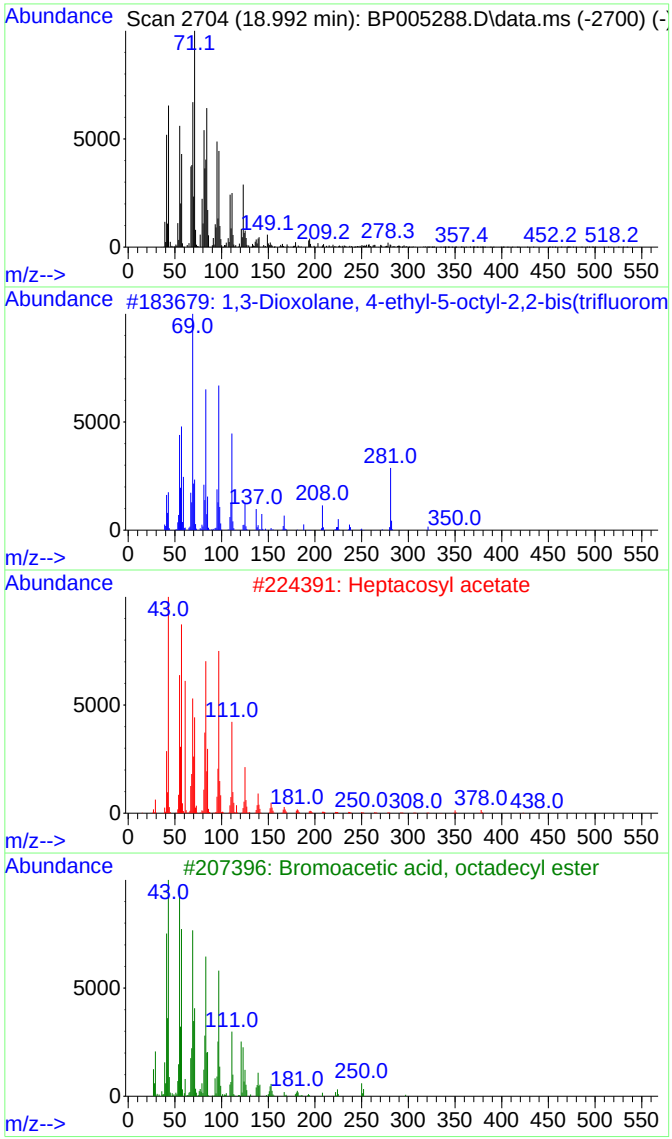
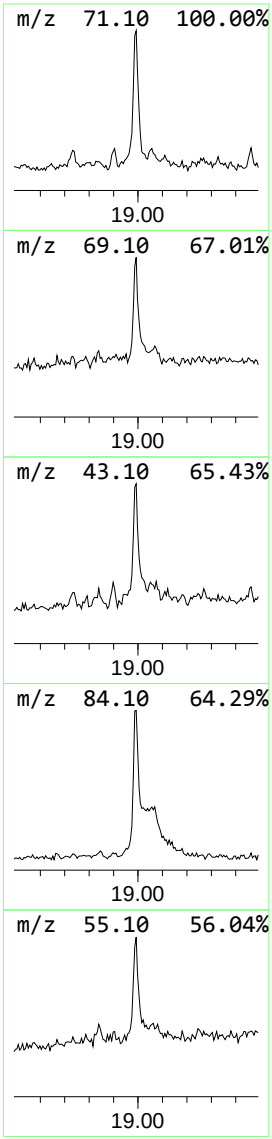
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1,3-Dioxolane, 4-ethyl-5-oc... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.992	9.75 ng	411784	Phenanthrene-d10	17.104

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Dioxolane, 4-ethyl-5-octyl-2...	350	C15H24F6O2	038274-73-6	55
2	Heptacosyl acetate	438	C29H58O2	1000351-78-2	46
3	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	38
4	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	023283-97-8	38
5	5-Eicosene, (E)-	280	C20H40	074685-30-6	35





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Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SB-5A

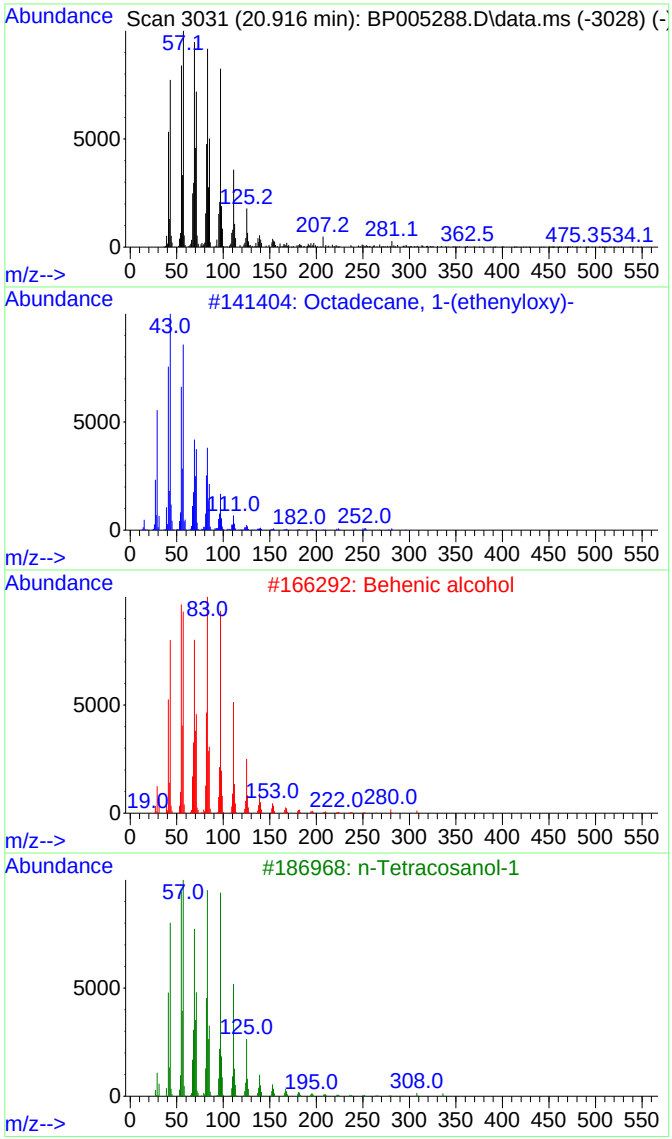
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

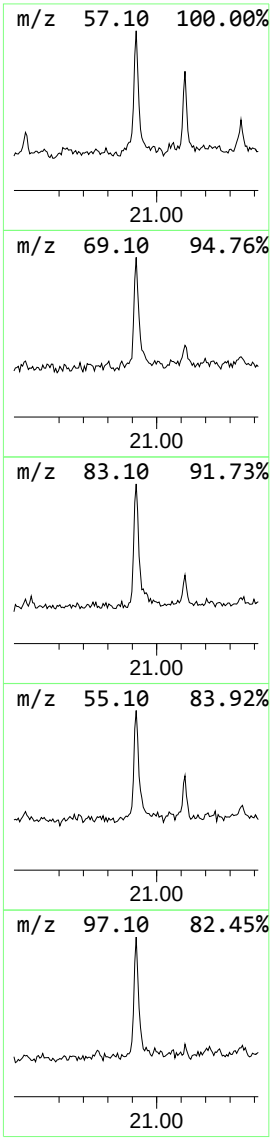
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Octadecane, 1-(ethenyloxy)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.916	10.55 ng	497626	Chrysene-d12	21.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	95
2	Behenic alcohol	326	C22H46O	000661-19-8	90
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	90
4	1-Heptacosanol	396	C27H56O	002004-39-9	90
5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	87





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Acq On : 13 Apr 2021 15:44
Operator : CG/JU
Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

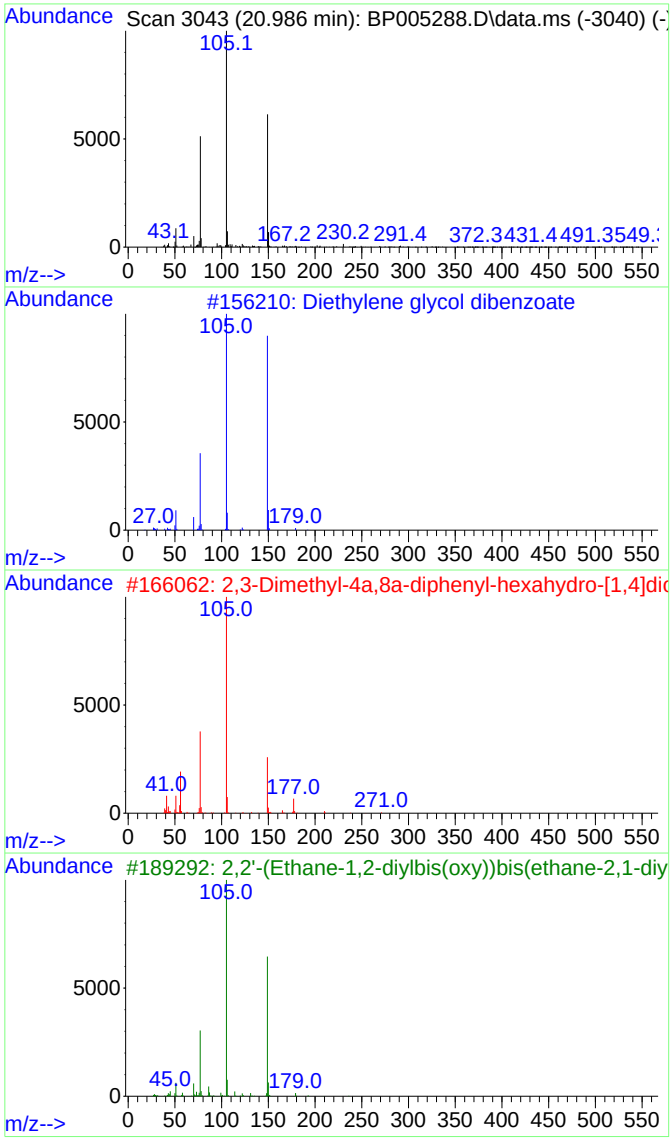
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

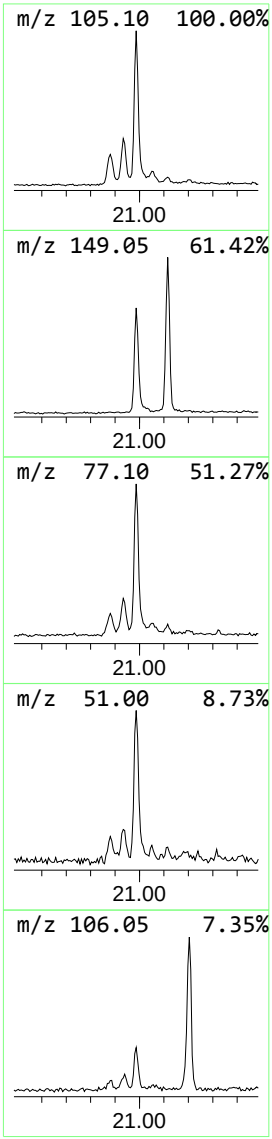
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Diethylene glycol dibenzoate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.986	5.40 ng	254678	Chrysene-d12	21.204

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	78
2	2,3-Dimethyl-4a,8a-diphenyl-hexa...	326	C20H22O4	1000297-07-0	53
3	2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	53
4	2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	47
5	Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	43





Tentatively Identified Compound (LSC) summary

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Sample : M1998-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : SB-5A

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1-Hexanol, 2-et...	7.740	9.6	ng	182634	1	7.675	380166	20.0
2-Piperidinone	10.299	52.8	ng	1284380	2	10.463	486590	20.0
Hexadecane, 2,6...	11.487	2.4	ng	58203	2	10.463	486590	20.0
Indole	11.987	36.9	ng	897120	2	10.463	486590	20.0
Pentadecane, 7-...	13.816	3.6	ng	126031	3	14.339	696114	20.0
2-Bromo dodecane	15.604	2.3	ng	80553	3	14.339	696114	20.0
Tridecane, 7-he...	16.086	6.4	ng	271430	4	17.104	844913	20.0
unknown16.875	16.875	9.1	ng	385761	4	17.104	844913	20.0
2-Pentadecanone...	17.198	4.0	ng	170216	4	17.104	844913	20.0
5,10-Diethoxy-2...	18.151	8.7	ng	366918	4	17.104	844913	20.0
1,3-Dioxolane, ...	18.992	9.8	ng	411784	4	17.104	844913	20.0
Octadecane, 1-(...	20.916	10.6	ng	497626	5	21.204	943393	20.0
Diethylene glyc...	20.986	5.4	ng	254678	5	21.204	943393	20.0