LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 18-B1(0-2)

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-BP052721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005667.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.117	3	5	10	rVB	2207688	2295758	27.13%	2.502%
2	3.158	10	12	21	rVB	175562	217403	2.57%	0.237%
3	5.081	332	339	350	rVB2	192642	299186	3.54%	0.326%
4	5.552	412	419	427	rBV	3336435	4928680	58.25%	5.371%
5	7.152	683	691	699	rBV	3262009	5085540	60.10%	5.542%
6	7.587	760	765	778	rVB	110968	206658	2.44%	0.225%
7	7.987	827	833	843	rBV	1168565	1942560	22.96%	2.117%
8	9.152	1024	1031	1038	rBV	2018577	3223382	38.09%	3.512%
9	10.622	1276	1281	1286	rBV	63898	121759	1.44%	0.133%
10	10.799	1300	1311	1316	rBV	1387811	2462230	29.10%	2.683%
11	10.846	1316	1319	1328	rVB	174580	277755	3.28%	0.303%
12	12.446	1587	1591	1599	rVB	62622	103412	1.22%	0.113%
13	13.240	1719	1726	1738	rBV	5077183	7400236	87.46%	8.064%
14	13.446	1756	1761	1767	rVB2	77049	112393	1.33%	0.122%
15	13.934	1838	1844	1849	rBV3	74145	132090	1.56%	0.144%
16	14.063	1859	1866	1870	rBV	508047	757999	8.96%	0.826%
17	14.334	1905	1912	1919	rBV	89012	189720	2.24%	0.207%
18	14.446	1927	1931	1938	rVB3	59582	109386	1.29%	0.119%
19	14.510	1938	1942	1951	rVB2	81850	115927	1.37%	0.126%
20	14.610	1951	1959	1965	rBV	1963276	2912015	34.41%	3.173%
21	14.675	1965	1970	1978	rVB	226917	354438	4.19%	0.386%
22	15.004	2021	2026	2033	rVB	203094	303195	3.58%	0.330%
23	15.457	2100	2103	2107	rVB	77292	92389	1.09%	0.101%
24	15.651	2131	2136	2148	rVB	253506	402909	4.76%	0.439%
25	15.863	2168	2172	2175	rBV2	73625	106723	1.26%	0.116%
26	15.945	2182	2186	2194	rVB	108062	199171	2.35%	0.217%
27	16.092	2205	2211	2219	rBV	3553229	5230054	61.81%	5.699%
28	16.322	2246	2250	2251	rBV	128274	154175	1.82%	0.168%
29	17.004	2362	2366	2371	rBV	126915	172388	2.04%	0.188%
30	17.110	2381	2384	2389	rVB2	127714	162540	1.92%	0.177%
31	17.169	2390	2394	2404	rVB2	279960	453230	5.36%	0.494%
32	17.351	2420	2425	2429	rBV	2145470	3209815	37.93%	3.498%
33	17.392	2429	2432	2439	rVV	3271985	4759378	56.25%	5.186%
34	17.487	2443	2448	2453	rVB	761972	1056035	12.48%	1.151%
35	17.751	2489	2493	2500	rBV	209992	301707	3.57%	0.329%
36	17.851	2507	2510	2517	rBV3	106429	214686	2.54%	0.234%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

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-BP052721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	17.922	2519	2522	2526	rVV2	129242	187967	2.22%	0.205%
38	18.222	2568	2573	2577	rBV2	570315	1074689	12.70%	1.171%
39	18.263	2577	2580	2586	rVB	515511	670026	7.92%	0.730%
40	18.339	2590	2593	2596	rVB	188071	196208	2.32%	0.214%
41	18.404	2599	2604	2607	rBV2	527301	836914	9.89%	0.912%
42	18.692	2650	2653	2657	rVB	288087	355719	4.20%	0.388%
43	18.734	2657	2660	2665	rBV2	182005	271222	3.21%	0.296%
44	19.092	2719	2721	2725	rBV	199950	254133	3.00%	0.277%
45	19.351	2760	2765	2771	rBV	5075854	6504694	76.87%	7.088%
46	19.704	2821	2825	2832	rVB	4157317	5434297	64.22%	5.922%
47	19.898	2853	2858	2862	rVB	6695111	8461508	100.00%	9.220%
48	20.328	2927	2931	2937	rBV2	1266452	1879302	22.21%	2.048%
49	21.416	3111	3116	3120	rBV2	3145632	6471226	76.48%	7.052%
50	21.463	3121	3124	3132	rVB	1997900	2863893	33.85%	3.121%
51	21.933	3201	3204	3208	rVB	1212051	1485775	17.56%	1.619%
52	23.122	3402	3406	3411	rBV	1325086	2502204	29.57%	2.727%
53	23.757	3511	3514	3520	rVB	1341467	2252375	26.62%	2.454%

Sum of corrected areas: 91769074

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Operator : CG/JU
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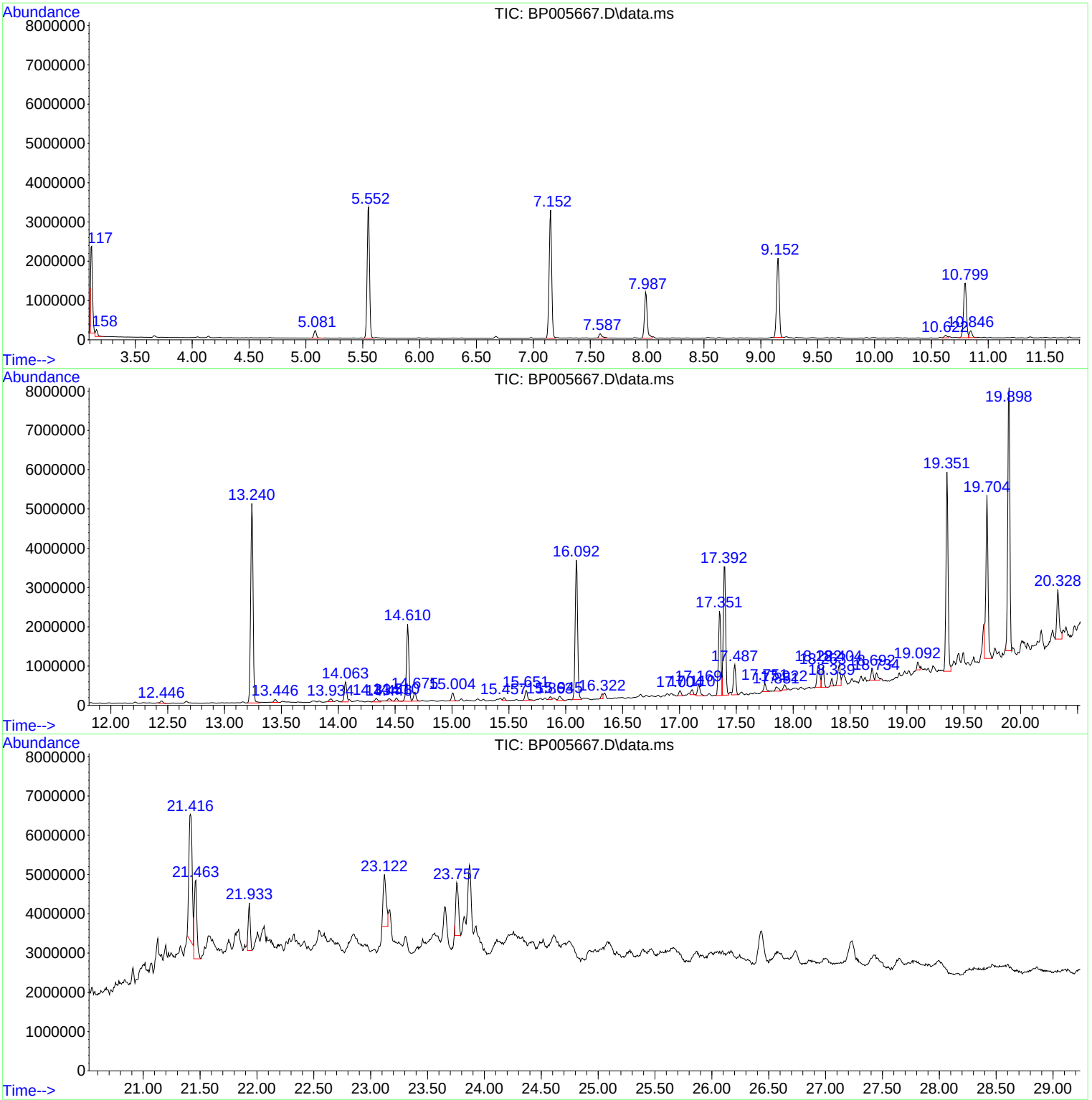
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

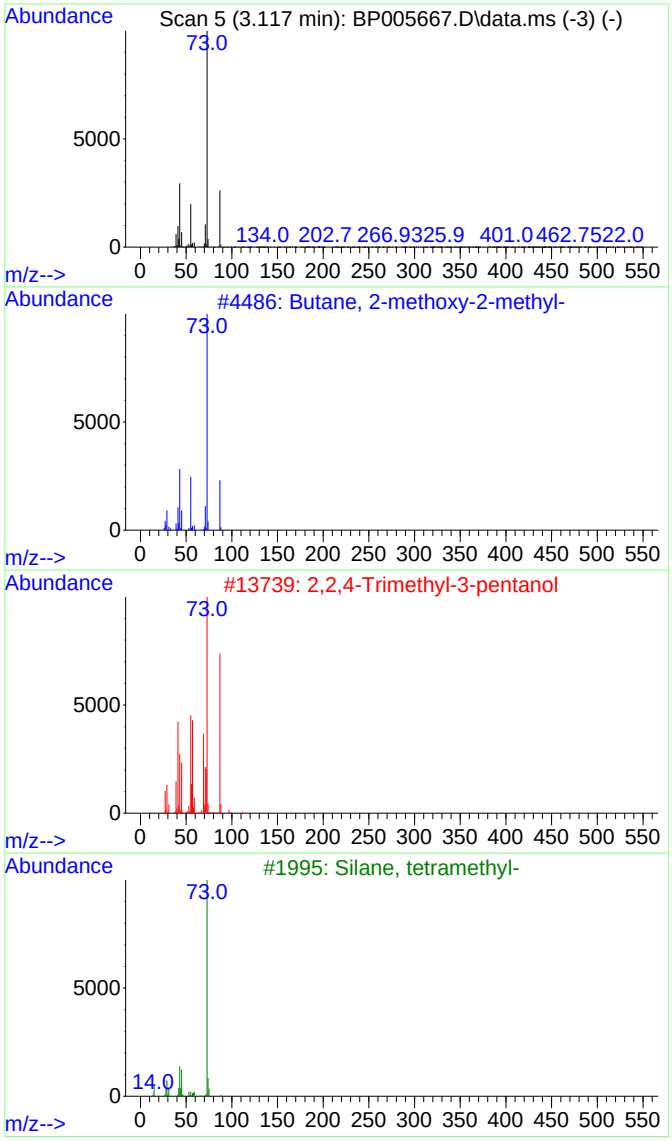
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

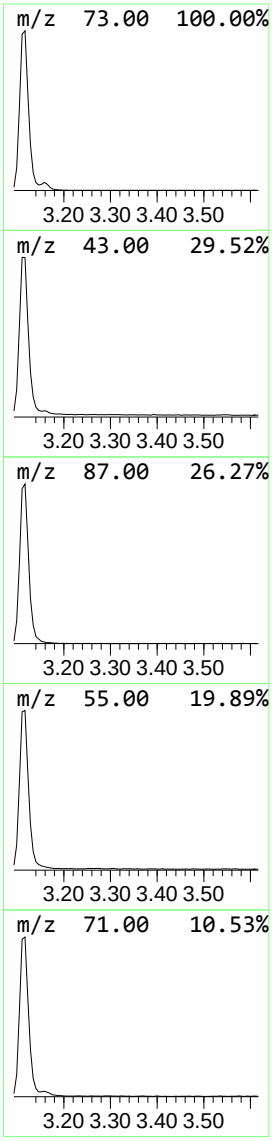
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.117	23.64 ng	2295760	1,4-Dichlorobenzene-d4	7.987

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

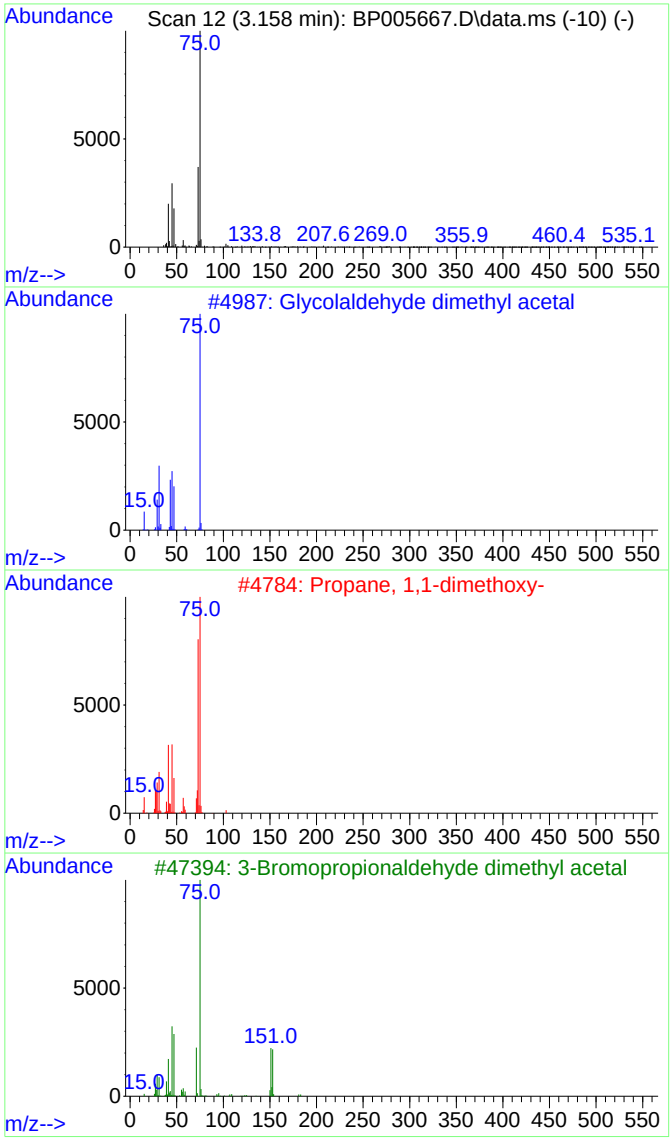
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

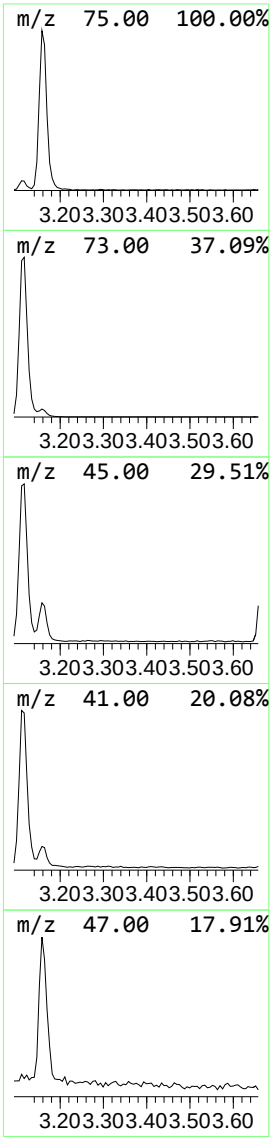
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Glycolaldehyde dimethyl acetal Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.158	2.24 ng	217403	1,4-Dichlorobenzene-d4	7.987

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	50
2	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	39
3	3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	9
4	Ethane, 1,1,2,2-tetramethoxy-	150	C6H14O4	002517-44-4	9
5	Benzene, [2-(ethylthio)ethyl]-	166	C10H14S	022914-08-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 18-B1(0-2)

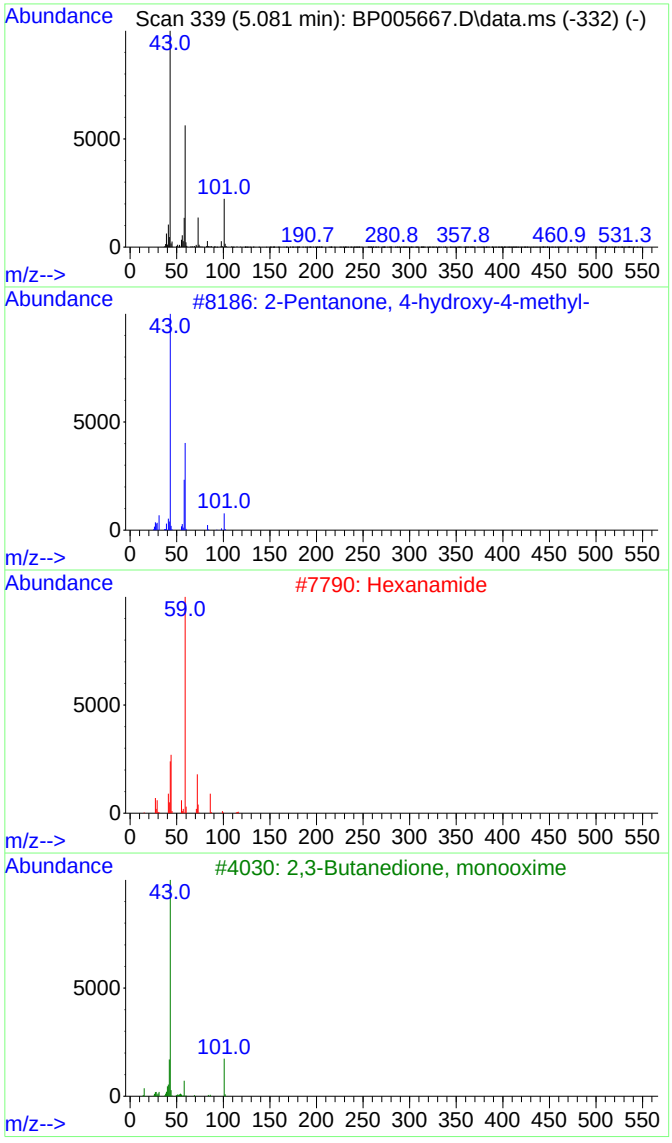
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

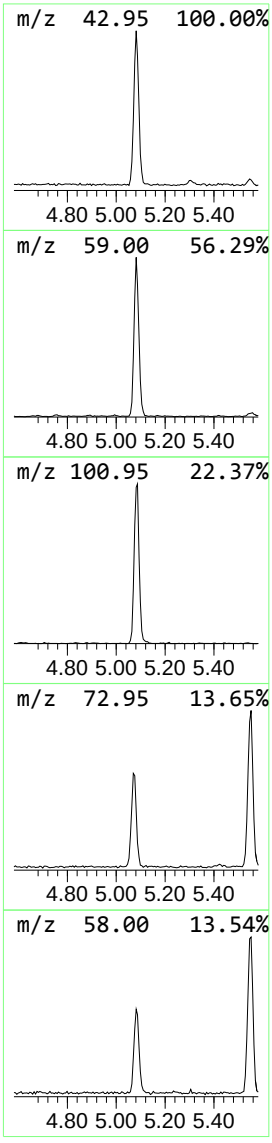
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.081	3.08 ng	299186	1,4-Dichlorobenzene-d4	7.987

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	Hexanamide	115	C6H13NO	000628-02-4	35
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	23
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
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Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

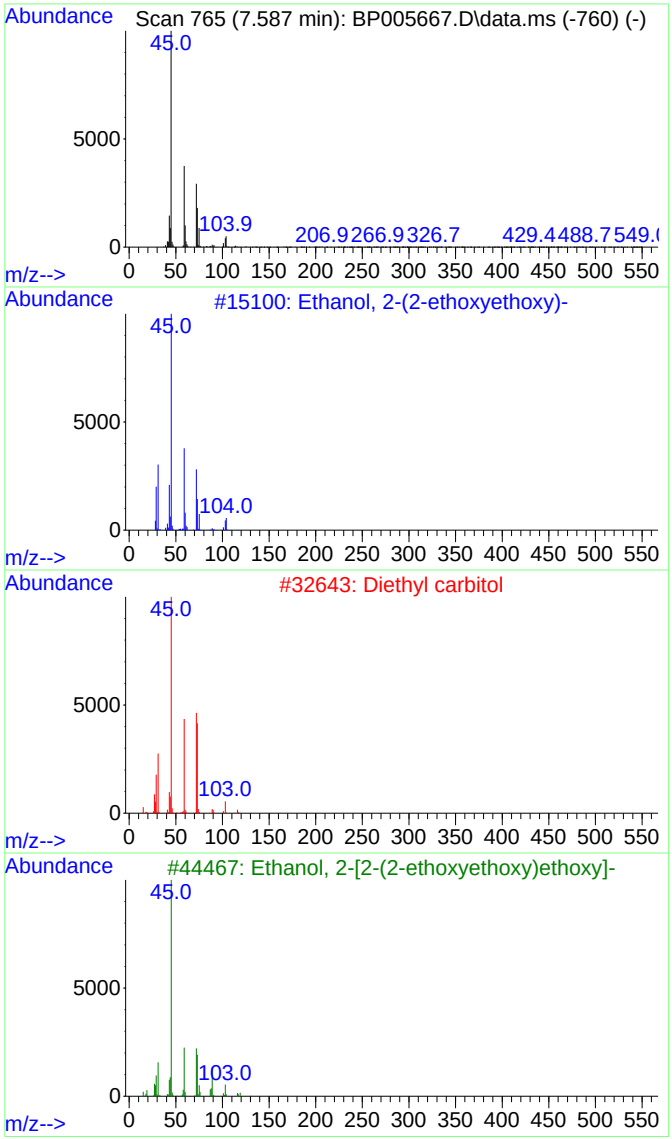
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

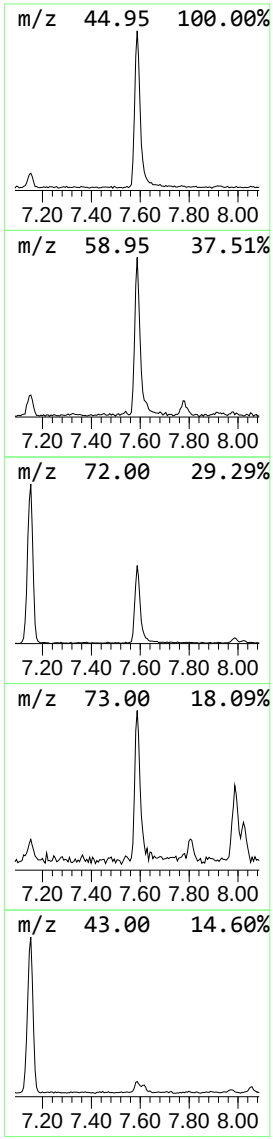
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.587	2.13 ng	206658	1,4-Dichlorobenzene-d4	7.987

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2	Diethyl carbitol	162	C8H18O3	000112-36-7	72
3	Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	56
4	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	45
5	Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
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Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

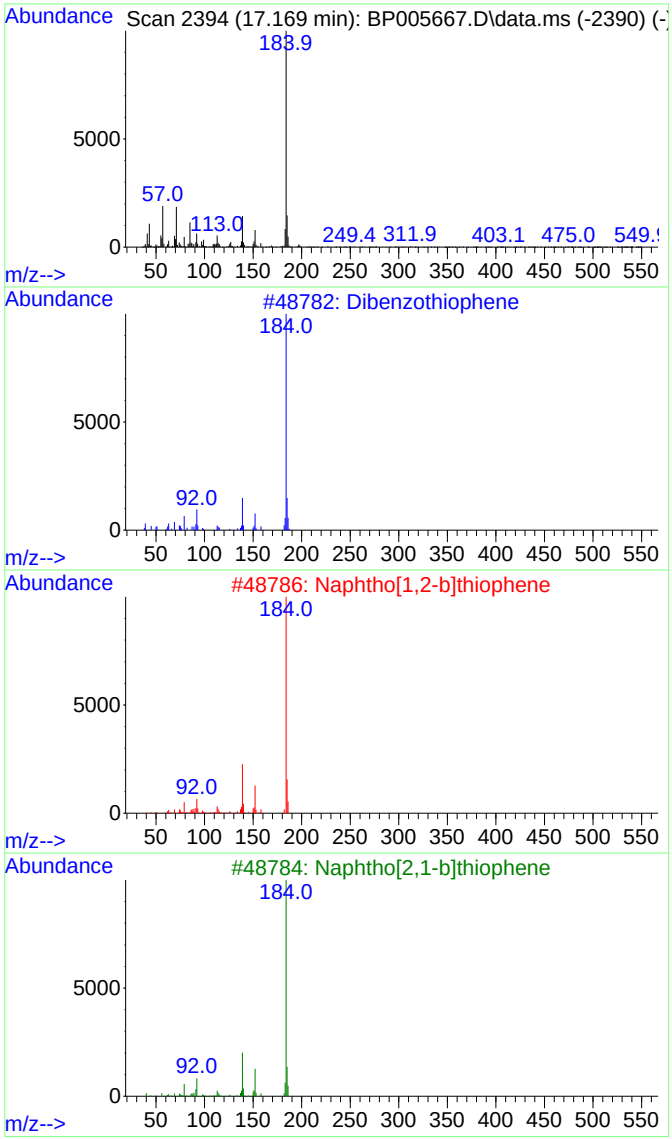
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

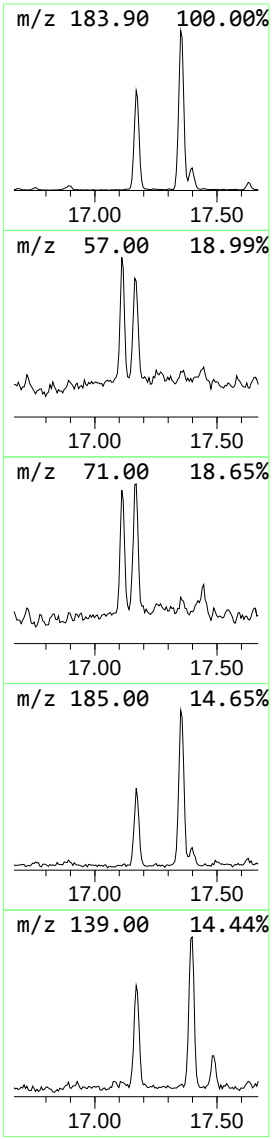
****

Peak Number 8 Dibenzothiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.169	2.82 ng	453230	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	91
2	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	87
3	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	87
4	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	87
5	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	87





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Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

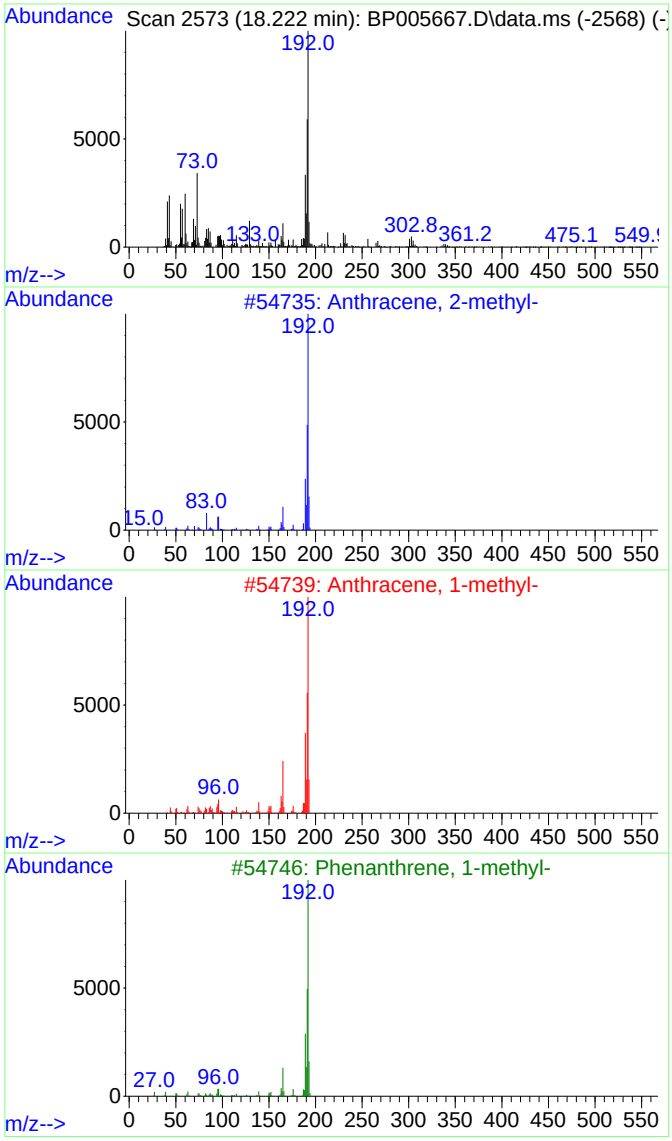
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

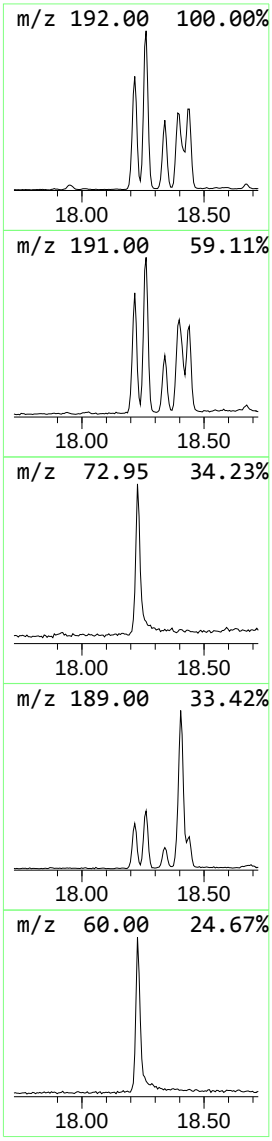
****

Peak Number 9 Anthracene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	6.70 ng	1074690	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	70
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	70





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Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

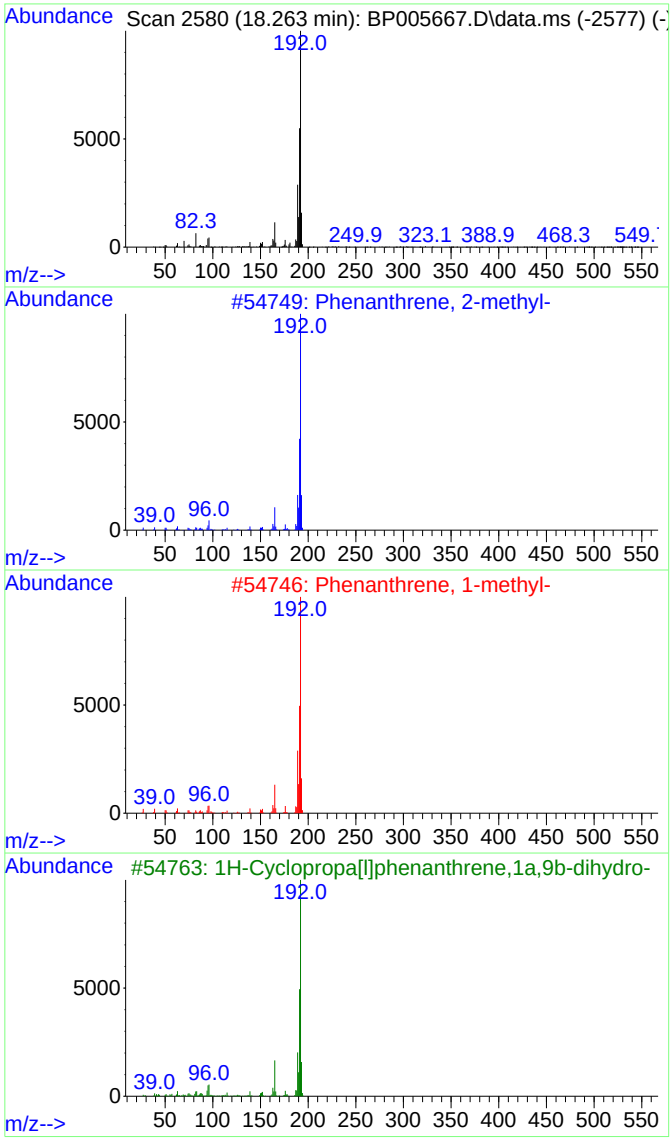
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

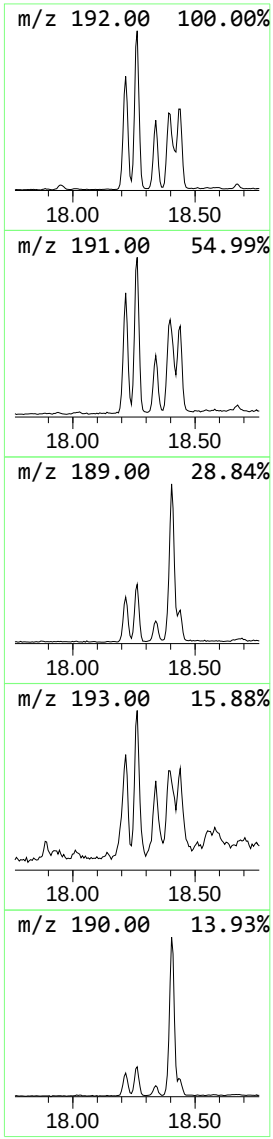
****

Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.263	4.17 ng	670026	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

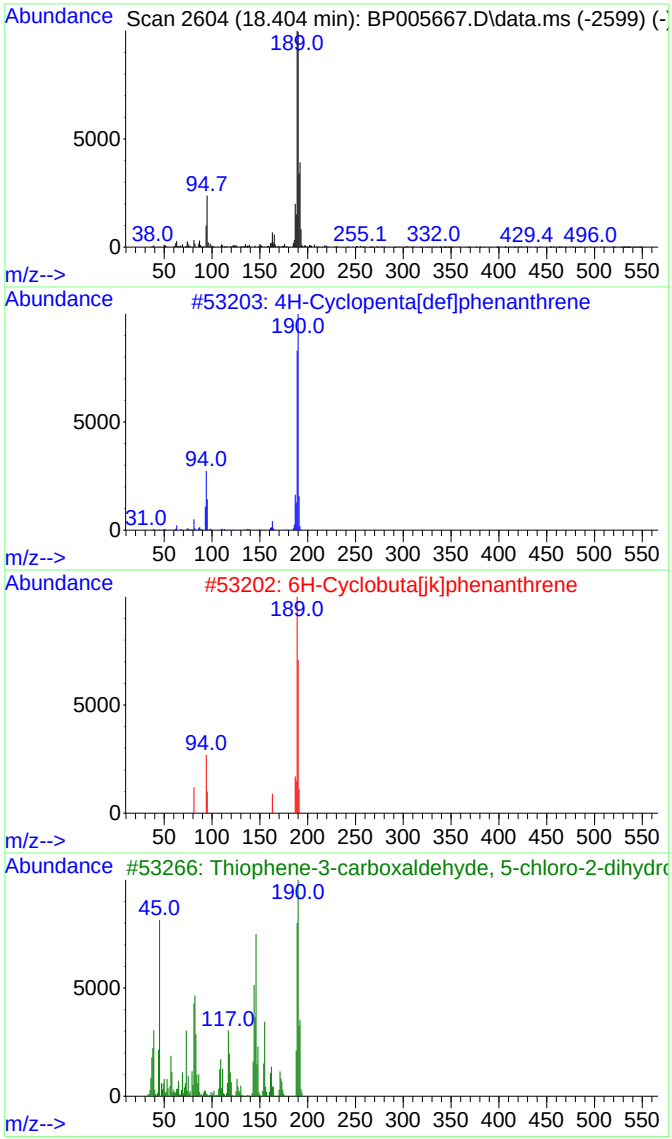
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

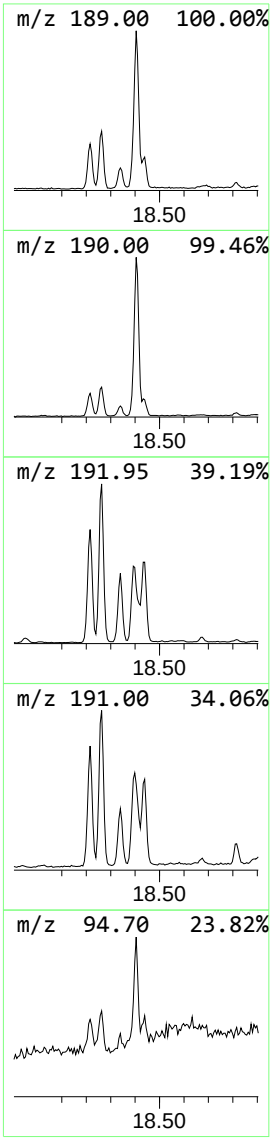
****

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.404	5.21 ng	836914	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	40
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38





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Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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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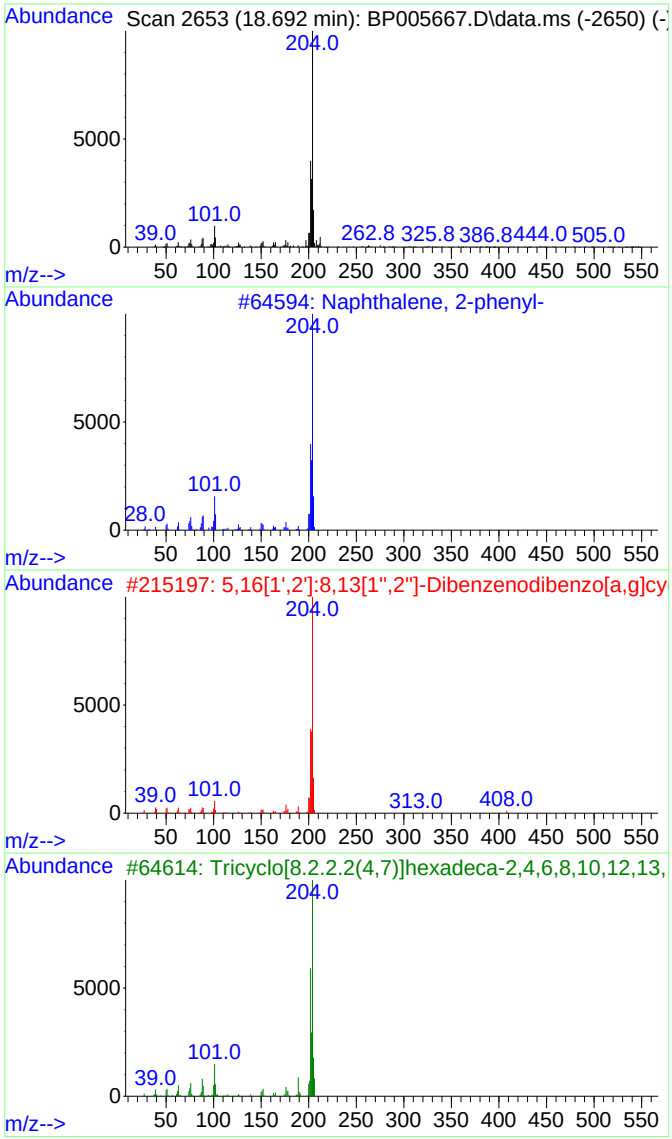
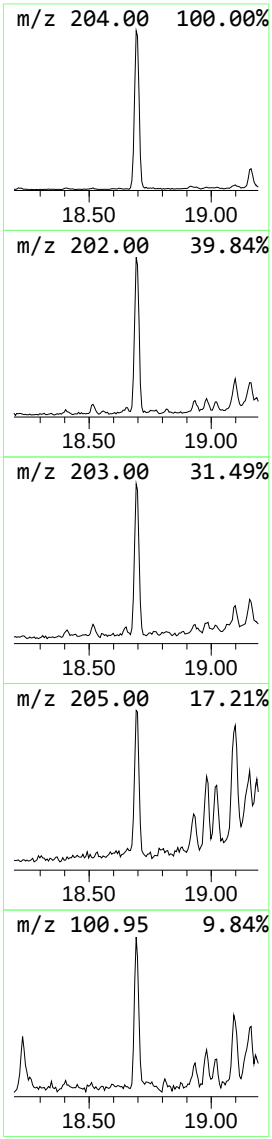
Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.692	2.22 ng	355719	Phenanthrene-d10	17.351

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	91
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	64
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
4	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	50
5	Levamisole	204	C11H12N2S	014769-73-4	49





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Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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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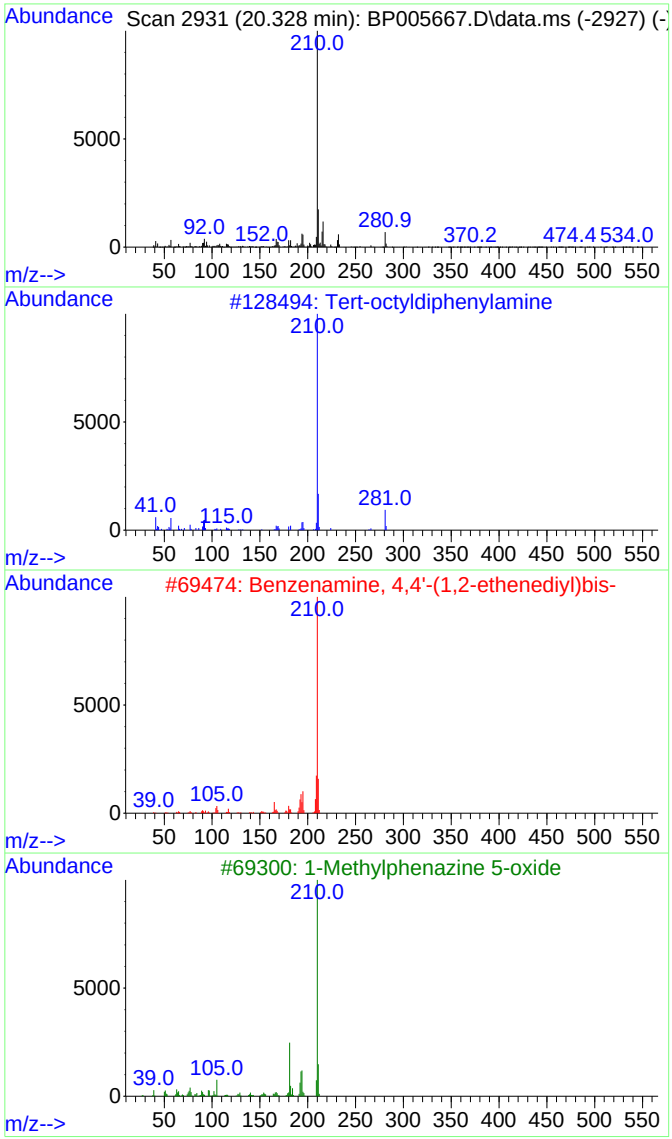
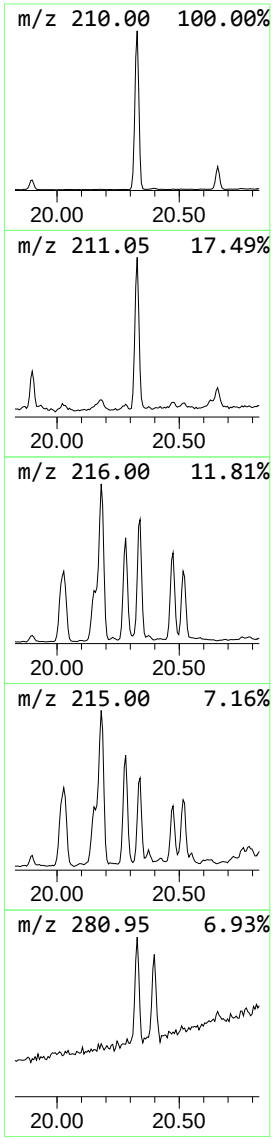
Peak Number 13 Tert-octyldiphenylamine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.328	5.81 ng	1879300	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tert-octyldiphenylamine	281	C20H27N	1000370-31-3	95
2	Benzenamine, 4,4'-(1,2-ethenediy...	210	C14H14N2	000621-96-5	59
3	1-Methylphenazine 5-oxide	210	C13H10N2O	014202-95-0	59
4	1-Acetoxy-6-methylphenazine	252	C15H12N2O2	014031-09-5	59
5	Cyclohepta[b]naphthalene-1-one	210	C15H14O	1000284-47-2	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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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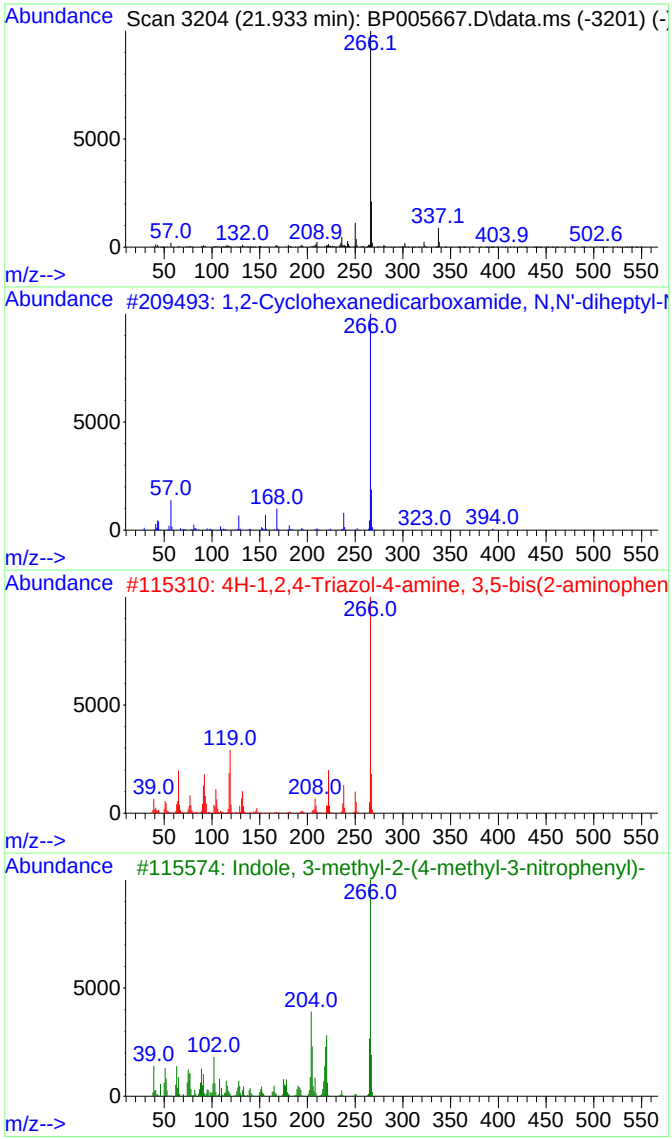
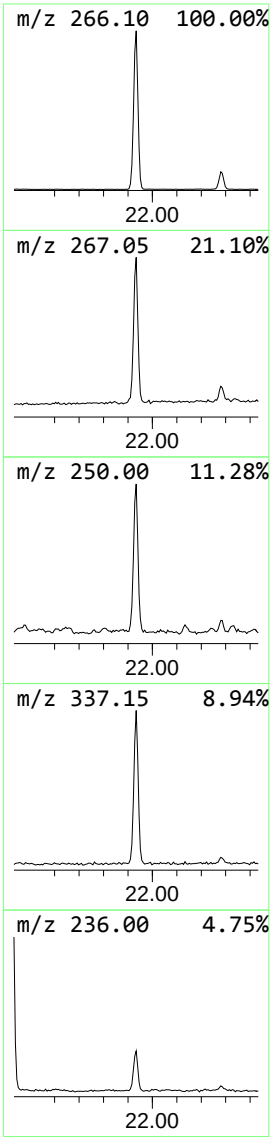
Peak Number 14 1,2-Cyclohexanedicarboxamid... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.933	4.59 ng	1485780	Chrysene-d12	21.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Cyclohexanedicarboxamide, N,...	394	C24H46N2O2	1000195-51-4	59
2	4H-1,2,4-Triazol-4-amine, 3,5-bi...	266	C14H14N6	1000350-00-5	56
3	Indole, 3-methyl-2-(4-methyl-3-n...	266	C16H14N2O2	311765-55-6	50
4	n-Pentafluorosulfanyl-S,S-diphen...	359	C12H10F5NO2S2	080039-31-2	50
5	Benzoic acid, 4-(4,5-dihydro-3-p...	266	C16H14N2O2	026192-74-5	45





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060221\
Data File : BP005667.D
Acq On : 02 Jun 2021 16:06
Operator : CG/JU
Sample : M2580-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : 18-B1(0-2)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.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
Butane, 2-metho...	3.117	23.6	ng	2295760	1	7.987	1942560	20.0
Glycolaldehyde ...	3.158	2.2	ng	217403	1	7.987	1942560	20.0
2-Pentanone, 4-...	5.081	3.1	ng	299186	1	7.987	1942560	20.0
Ethanol, 2-(2-e...	7.587	2.1	ng	206658	1	7.987	1942560	20.0
Dibenzothiophene	17.169	2.8	ng	453230	4	17.351	3209820	20.0
Anthracene, 2-m...	18.222	6.7	ng	1074690	4	17.351	3209820	20.0
Phenanthrene, 2...	18.263	4.2	ng	670026	4	17.351	3209820	20.0
4H-Cyclopenta[d...	18.404	5.2	ng	836914	4	17.351	3209820	20.0
Naphthalene, 2-...	18.692	2.2	ng	355719	4	17.351	3209820	20.0
Tert-octyldiphe...	20.328	5.8	ng	1879300	5	21.422	6471230	20.0
1,2-Cyclohexane...	21.933	4.6	ng	1485780	5	21.422	6471230	20.0