LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

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\8270-BP091619.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.405	560	564	567	rBV	3426794	2877797	67.48%	4.855%
2	6.416	732	736	743	rVB	3693806	3174791	74.44%	5.356%
3	6.558	756	760	763	rBV	3986467	3244136	76.07%	5.473%
4	6.787	796	799	803	rBV	2401147	1889644	44.31%	3.188%
5	6.946	822	826	829	rVB	3480111	2705561	63.44%	4.564%
6	7.346	887	894	897	rBV	2212436	1851971	43.42%	3.124%
7	7.940	992	995	999	rBV	56843	54876	1.29%	0.093%
8	8.046	1009	1013	1014	rBV	78134	60824	1.43%	0.103%
9	8.069	1014	1017	1020	rVV	3102601	2419989	56.74%	4.082%
10	8.093	1020	1021	1028	rVB	154734	81755	1.92%	0.138%
11	8.499	1087	1090	1095	rBV	72988	74183	1.74%	0.125%
12	8.781	1134	1138	1141	rVB	56148	54666	1.28%	0.092%
13	9.146	1197	1200	1204	rBV	5292415	4264794	100.00%	7.195%
14	9.240	1212	1216	1219	rBV3	70957	85836	2.01%	0.145%
15	9.540	1264	1267	1273	rBV	487127	401300	9.41%	0.677%
16	9.687	1288	1292	1296	rBV	212299	188877	4.43%	0.319%
17	9.769	1304	1306	1311	rVB	66406	53624	1.26%	0.090%
18	9.828	1313	1316	1320	rVV	3422137	2802474	65.71%	4.728%
19	9.863	1320	1322	1325	rVB	72956	58464	1.37%	0.099%
20	10.034	1348	1351	1354	rBV	91203	71735	1.68%	0.121%
21	10.381	1407	1410	1416	rBV2	80542	114466	2.68%	0.193%
22	10.552	1435	1439	1442	rBV2	120858	126696	2.97%	0.214%
23	10.593	1442	1446	1447	rBV	153340	142131	3.33%	0.240%
24	10.616	1447	1450	1456	rVB	2646312	2461464	57.71%	4.152%
25	10.757	1469	1474	1477	rBV	730774	683459	16.03%	1.153%
26	10.810	1480	1483	1485	rVV	1611919	1274966	29.89%	2.151%
27	10.846	1485	1489	1492	rVV2	1438804	1772220	41.55%	2.990%
28	10.875	1492	1494	1497	rVV2	1629644	1451579	34.04%	2.449%
29	10.899	1497	1498	1501	rVV	900432	692048	16.23%	1.167%
30	10.946	1501	1506	1507	rVV2	569825	779919	18.29%	1.316%
31	10.987	1510	1513	1517	rVV3	1188024	1383474	32.44%	2.334%
32	11.028	1517	1520	1522	rVV	1599661	1359249	31.87%	2.293%
33	11.046	1522	1523	1525	rVV	422750	360094	8.44%	0.607%
34	11.069	1525	1527	1531	rVB2	928132	716634	16.80%	1.209%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

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\8270-BP091619.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.322	1567	1570	1573	rBV	3358319	2854574	66.93%	4.816%
36	11.346	1573	1574	1577	rVV	598063	343478	8.05%	0.579%
37	11.399	1580	1583	1586	rVB	323946	259561	6.09%	0.438%
38	11.434	1586	1589	1593	rBV4	47298	60868	1.43%	0.103%
39	11.551	1607	1609	1613	rBV	58869	56315	1.32%	0.095%
40	11.828	1654	1656	1659	rBV	90706	78499	1.84%	0.132%
41	11.857	1659	1661	1666	rBV	428018	430901	10.10%	0.727%
42	11.904	1667	1669	1672	rVB	83763	63861	1.50%	0.108%
43	11.940	1672	1675	1678	rBV	191288	229700	5.39%	0.387%
44	12.128	1705	1707	1709	rBV	77468	63961	1.50%	0.108%
45	12.387	1748	1751	1754	rBV2	112622	98754	2.32%	0.167%
46	12.546	1775	1778	1781	rBV	1492829	1184399	27.77%	1.998%
47	12.593	1784	1786	1792	rVV2	193248	192530	4.51%	0.325%
48	12.775	1812	1817	1819	rBV	1349674	1268401	29.74%	2.140%
49	12.804	1819	1822	1825	rVB	1728306	1593190	37.36%	2.688%
50	12.922	1839	1842	1846	rBV	4817910	4264893	100.00%	7.195%
51	13.110	1872	1874	1877	rVB	233028	196686	4.61%	0.332%
52	13.845	1996	1999	2007	rVB4	418442	565495	13.26%	0.954%
53	13.987	2019	2023	2026	rBV2	2992316	3458170	81.08%	5.834%
54	14.010	2026	2027	2030	rVB	614574	479989	11.25%	0.810%
55	15.475	2273	2276	2280	rVB	1594365	1797795	42.15%	3.033%

Sum of corrected areas: 59277716

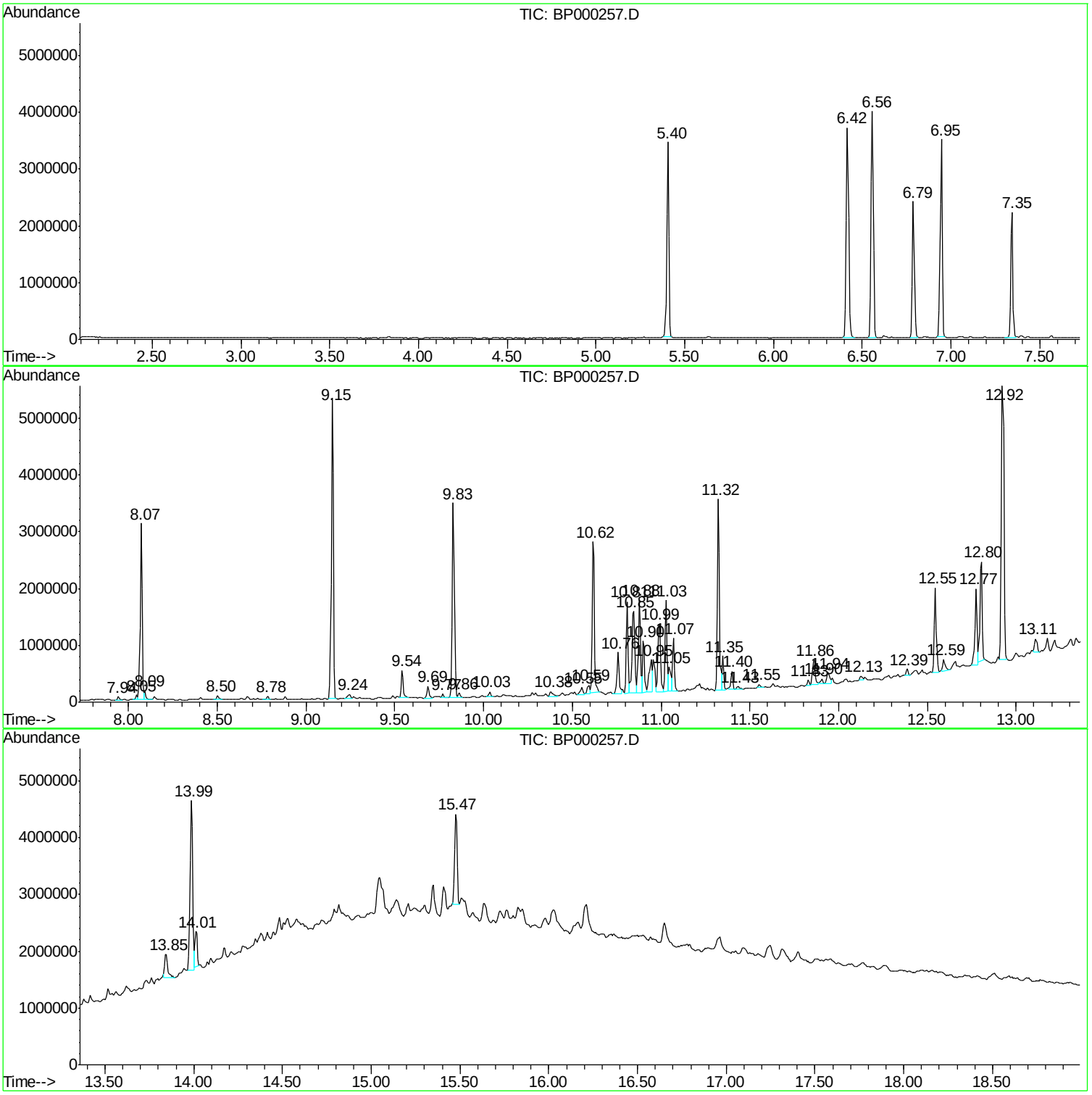
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

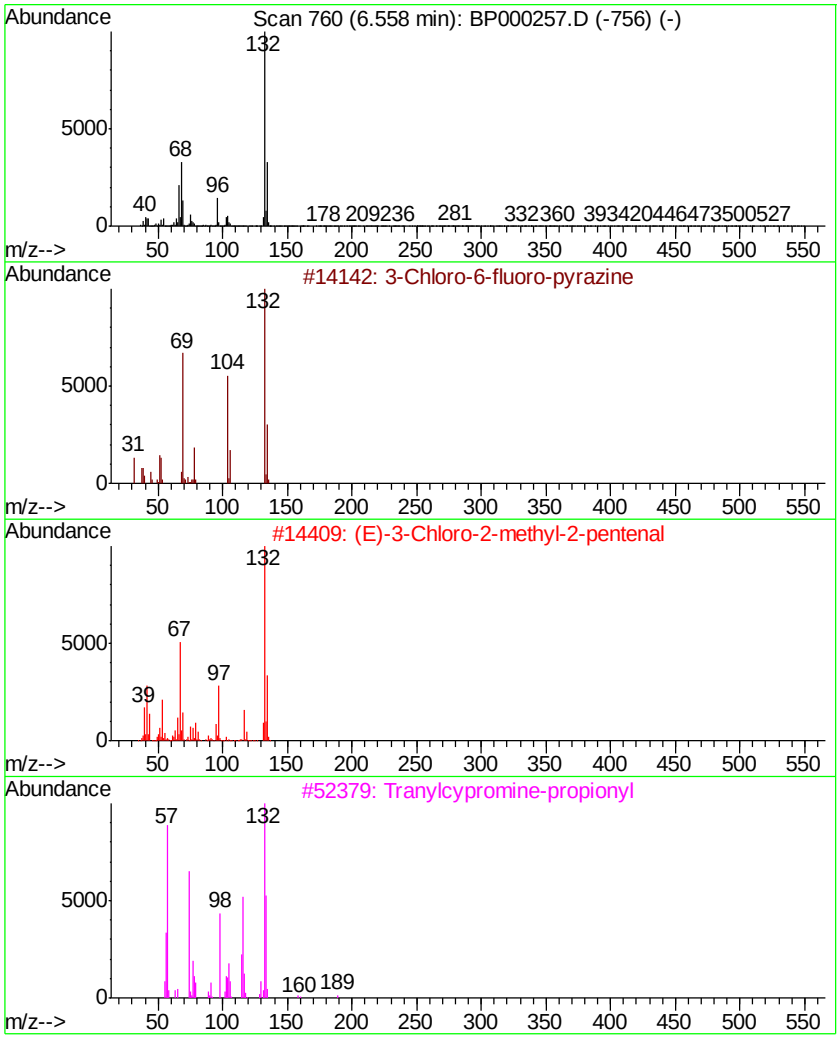
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

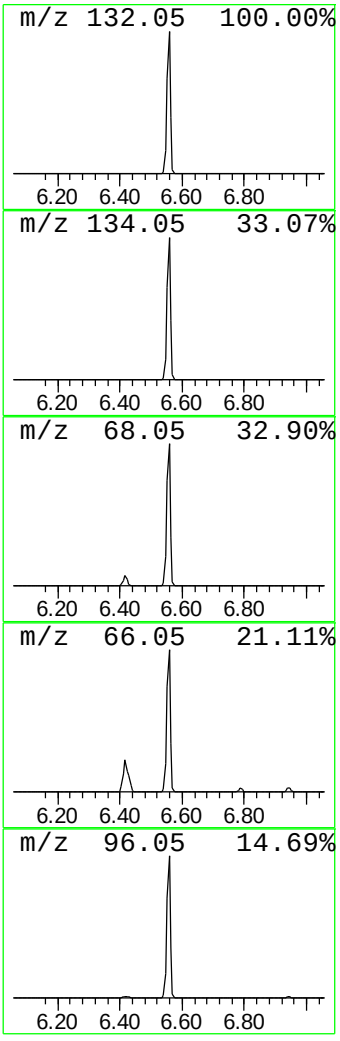
****

Peak Number 1 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	34.34 ng	3244140	1,4-Dichlorobenzene-d4	6.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

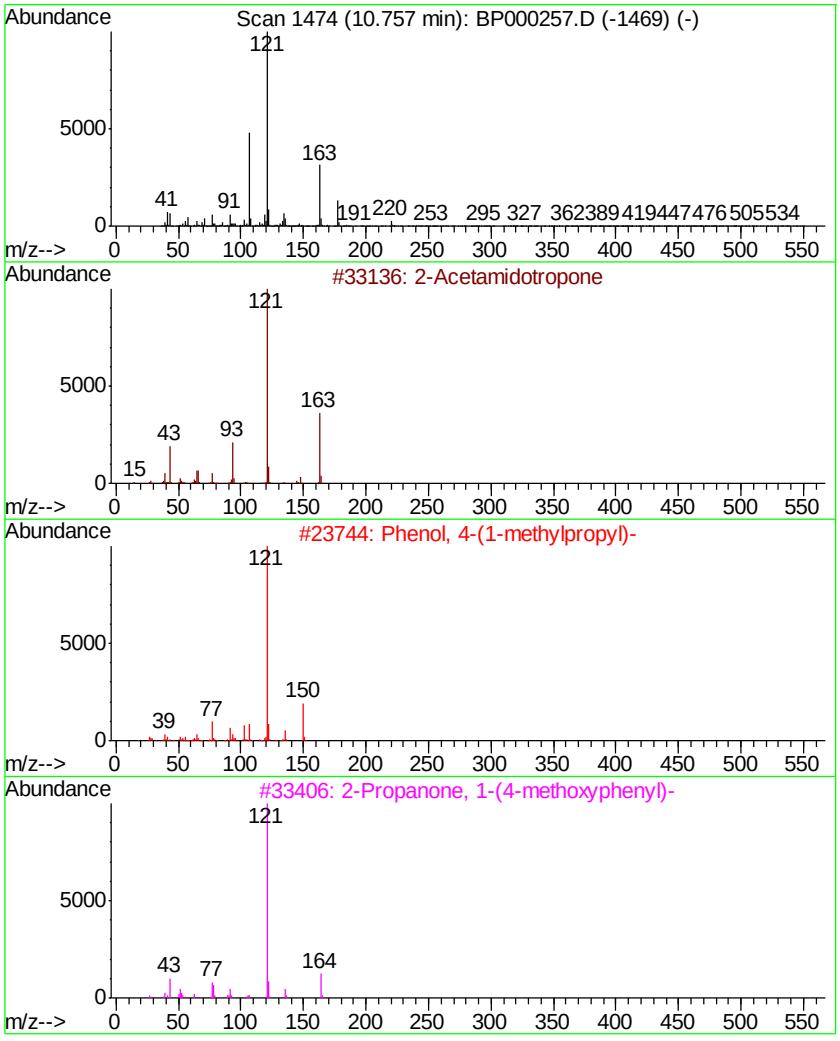
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

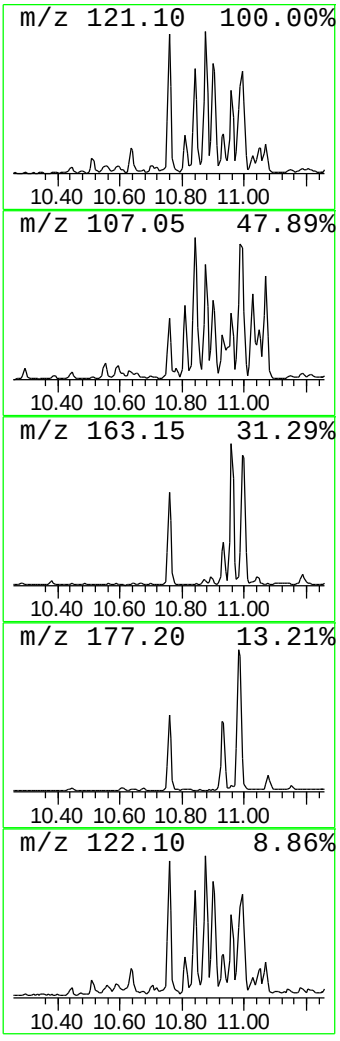
****

Peak Number 3 unknown10.76 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	4.79 ng	683459	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Acetamidotropone	163	C9H9NO2	006422-12-4	47
2	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	43
3	2-Propanone, 1-(4-methoxyphenyl)-	164	C10H12O2	000122-84-9	43
4	p-Sec-butylphenyl acetate	192	C12H16O2	003245-24-7	43
5	Acetamide, N-(2,6-dimethylphenyl)-	163	C10H13NO	002198-53-0	40





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

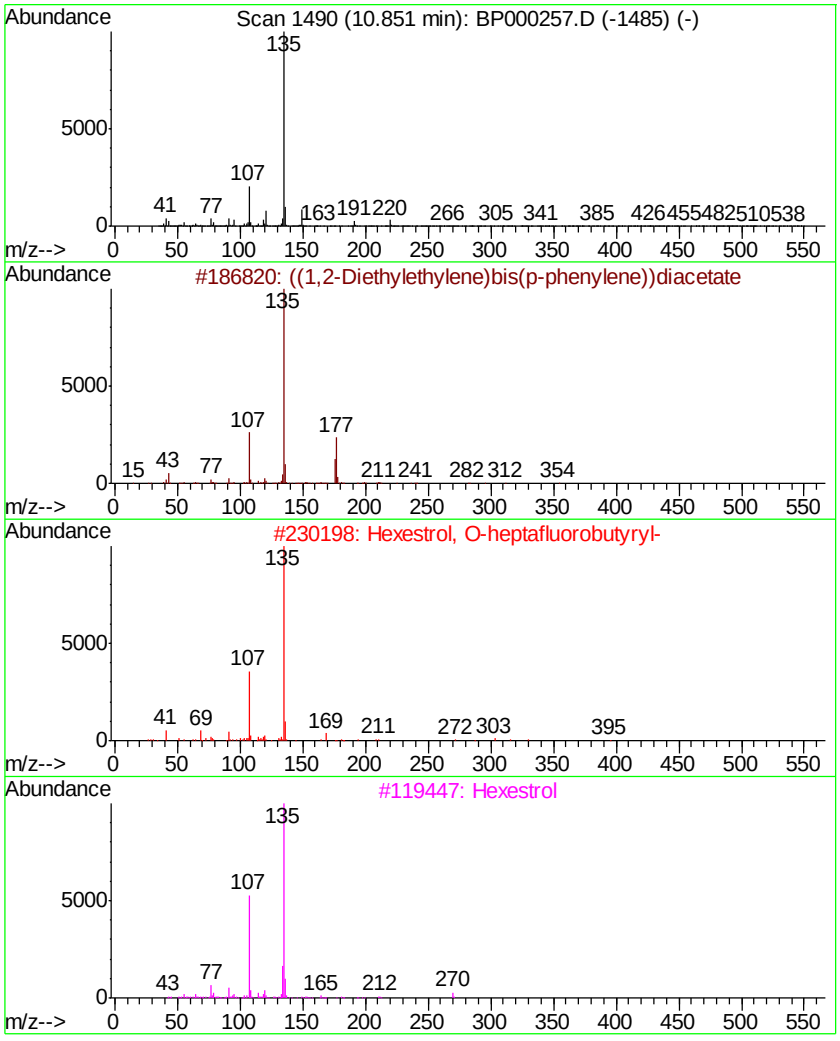
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

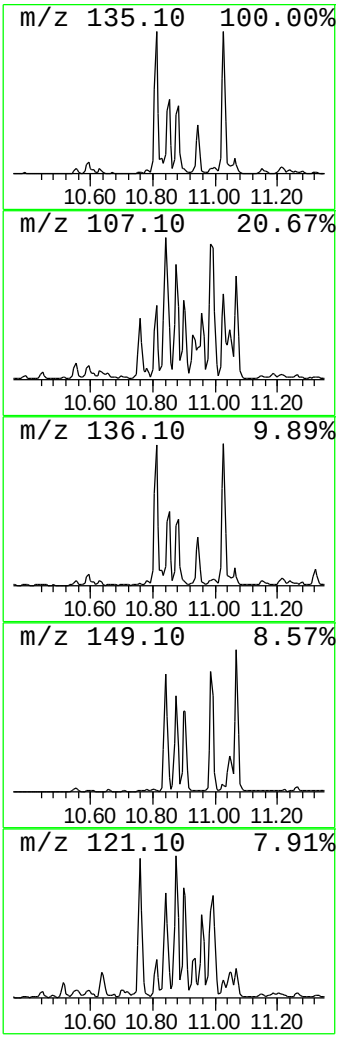
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 ((1,2-Diethylethylene)bis(p... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	12.42 ng	1772220	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	52
2	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	50
3	Hexestrol	270	C18H22O2	000084-16-2	50
4	Hexestrol	270	C18H22O2	005635-50-7	50
5	Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	50





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
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Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

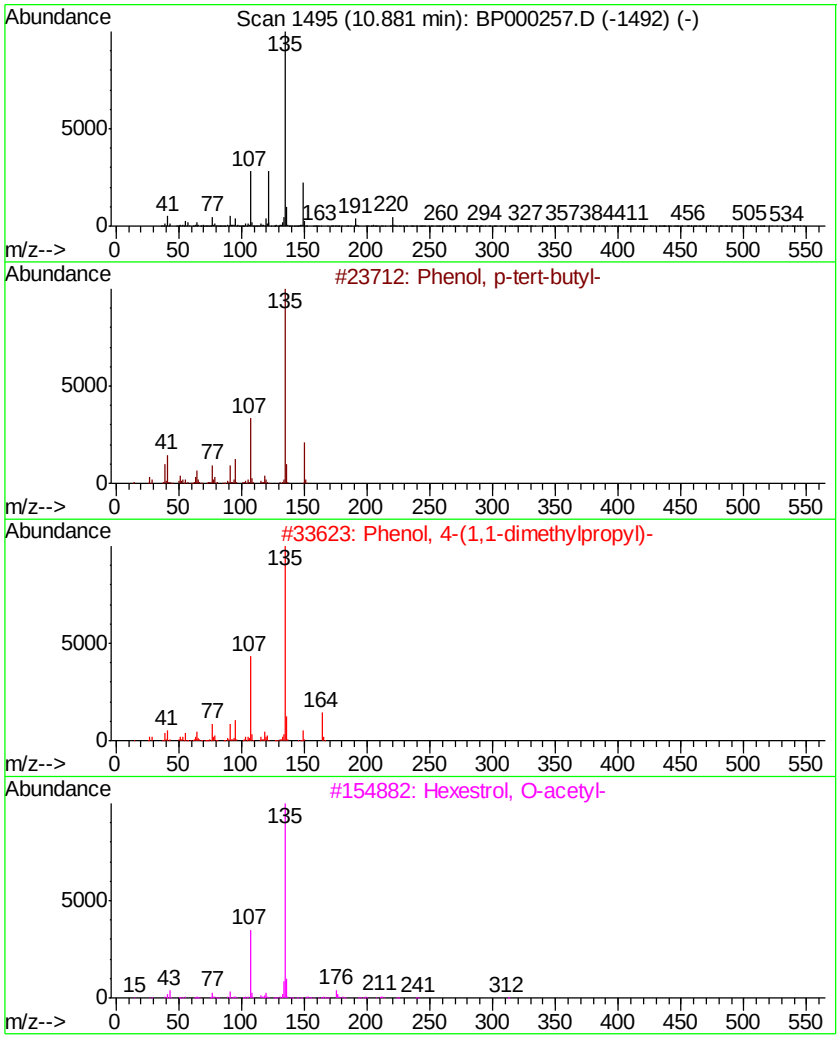
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

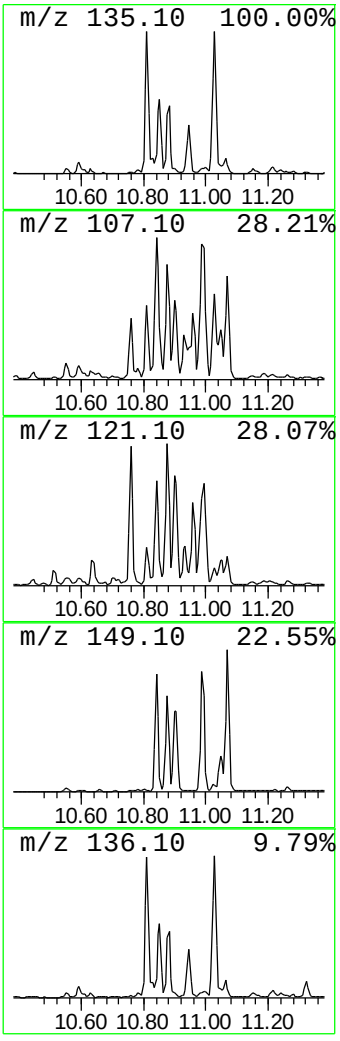
****

Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	10.17 ng	1451580	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	58
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50
3	Hexestrol, O-acetyl-	312	C20H24O3	1000374-87-8	50
4	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	50
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	50





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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

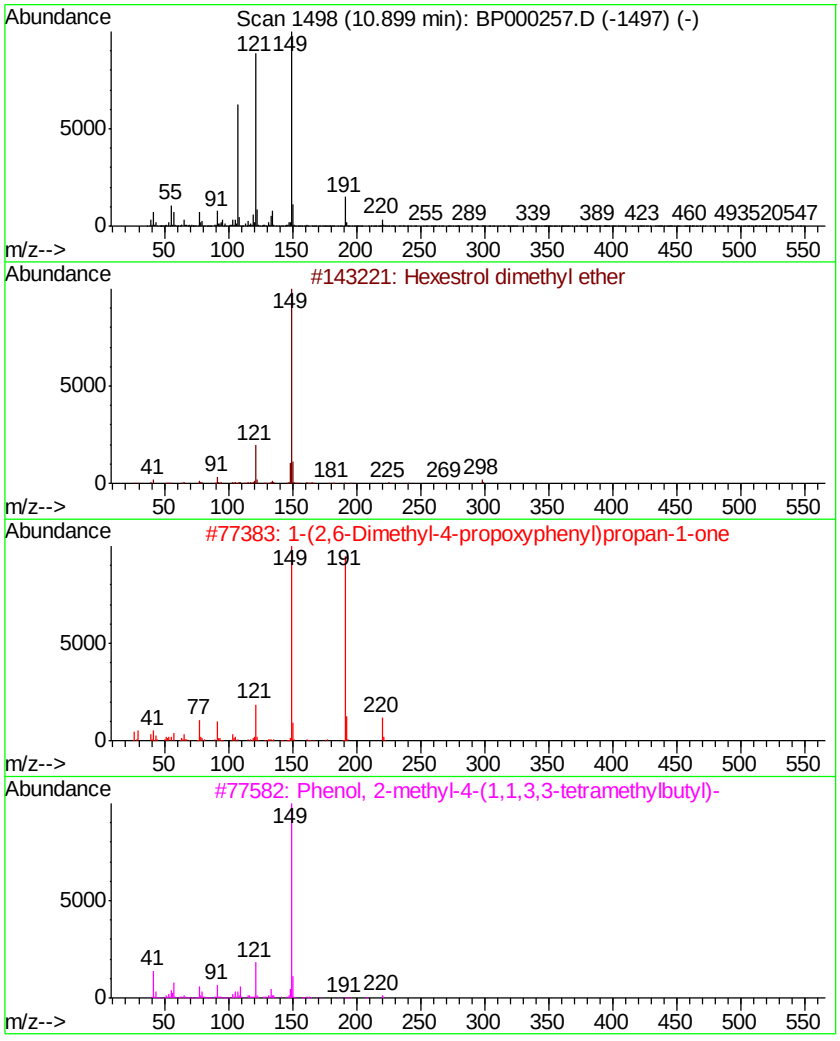
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

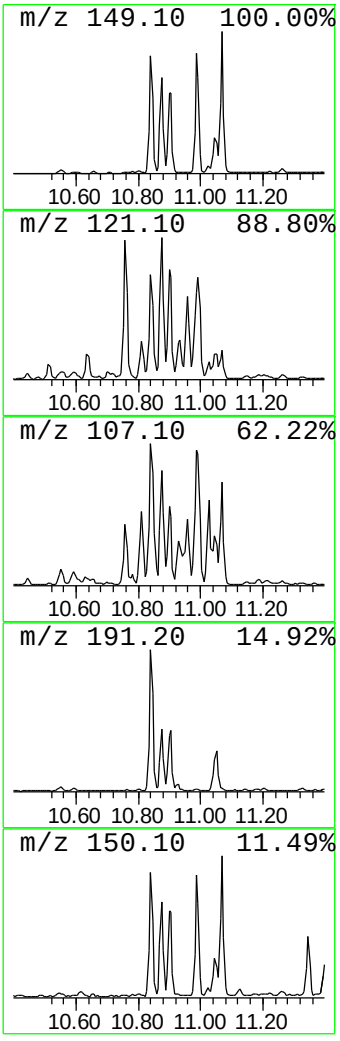
****

Peak Number 7 Hexestrol dimethyl ether Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	4.85 ng	692048	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexestrol dimethyl ether	298	C20H26O2	000130-78-9	64
2	1-(2,6-Dimethyl-4-propoxyphenyl)...	220	C14H20O2	1000190-22-3	35
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	35
4	Isoxazole, 4,5-dihydro-5-[(4-met...	191	C11H13NO2	1000362-14-0	30
5	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	30





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Misc :
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Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

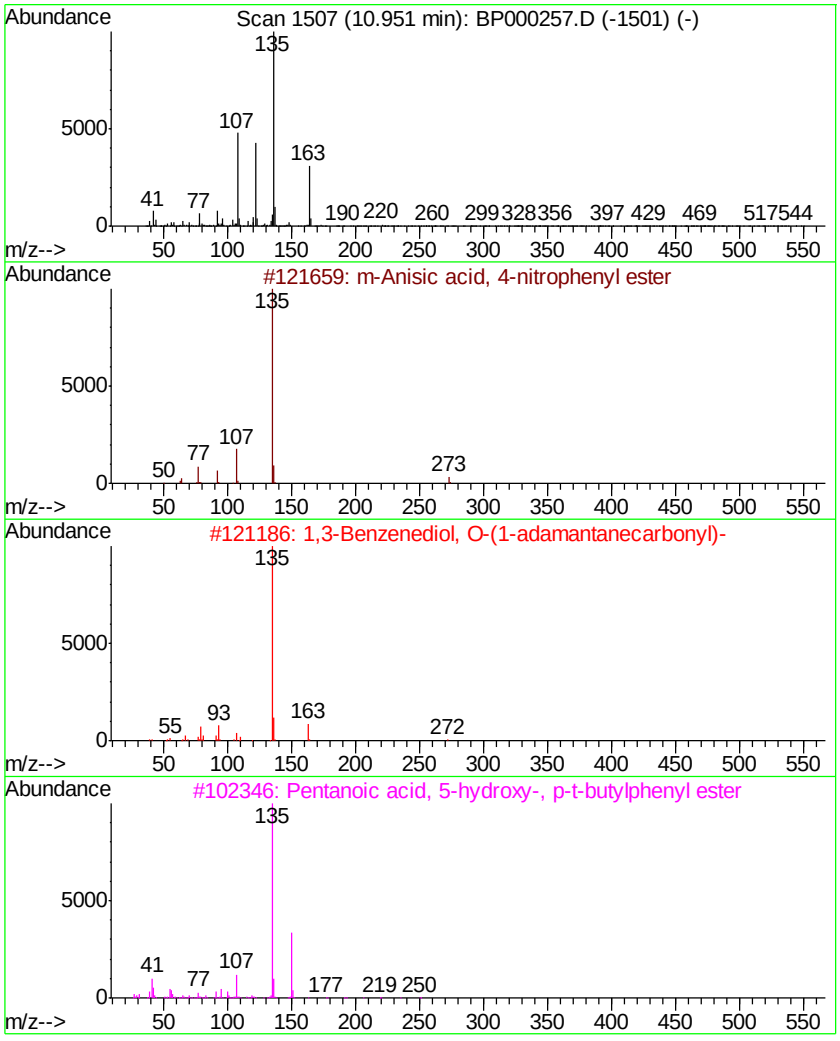
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

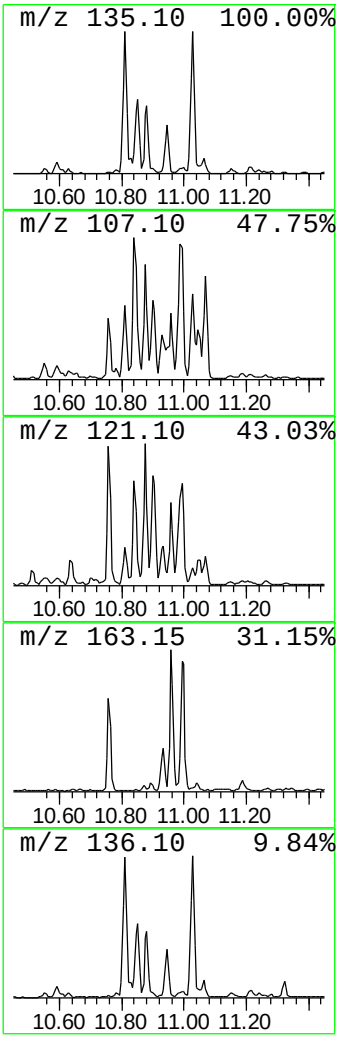
****

Peak Number 8 m-Anisic acid, 4-nitropheny... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.95	5.46 ng	779919	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Anisic acid, 4-nitrophenyl ester	273	C14H11NO5	1000307-80-4	72
2	1,3-Benzenediol, O-(1-adamantane...	272	C17H20O3	1000345-57-3	64
3	Pentanoic acid, 5-hydroxy-, p-t-...	250	C15H22O3	166273-37-6	64
4	4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	64
5	1,2-Benzenediol, o-(4-methoxyben...	362	C22H18O5	1000325-99-0	59





Library Search Compound Report

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Acq On : 16 Sep 2019 23:37
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Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

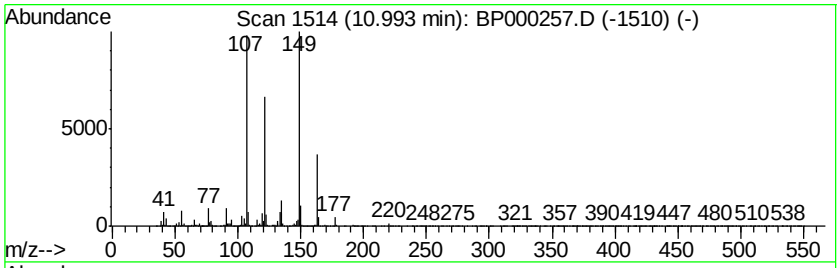
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

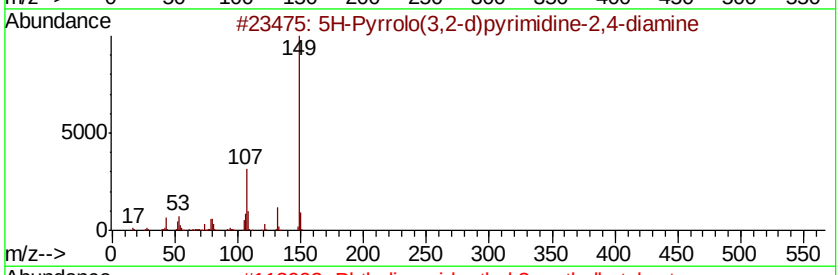
****

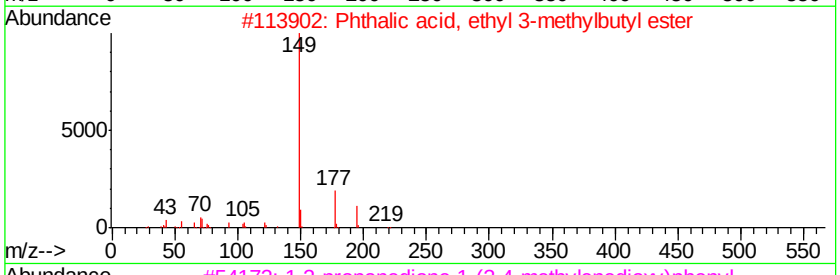
Peak Number 9 unknown10.99 Concentration Rank 5

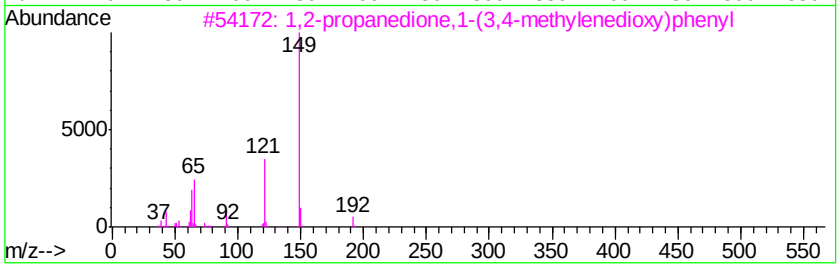
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	9.69 ng	1383470	Phenanthrene-d10	11.32

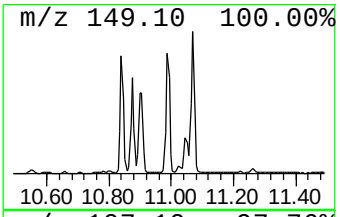
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	38
2	Phthalic acid, ethyl 3-methylbut...	264	C15H20O4	1000356-80-1	35
3	1,2-propanedione,1-(3,4-methylen...	192	C10H8O4	1000378-93-0	35
4	Phthalic acid, 2,7-dimethyloct-7...	370	C23H30O4	1000315-49-1	35
5	Hexestrol dimethyl ether	298	C20H26O2	000130-78-9	35

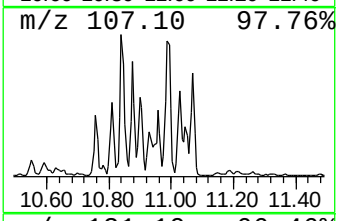


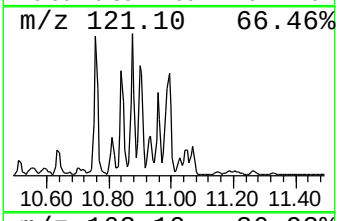


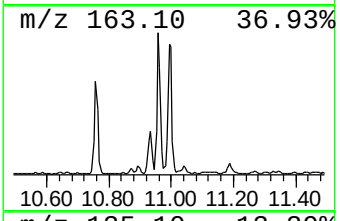


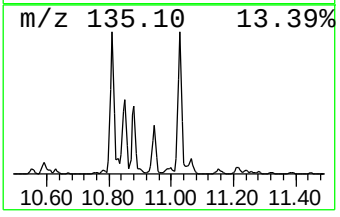












Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

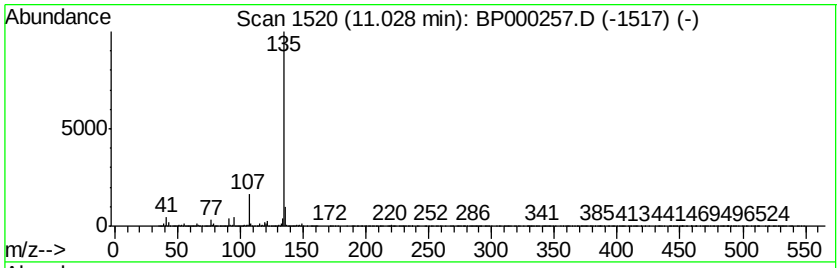
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

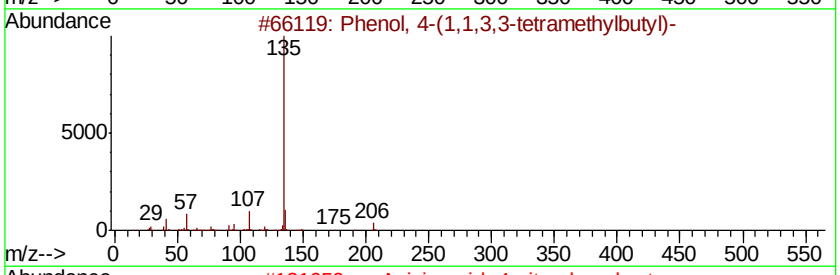
****

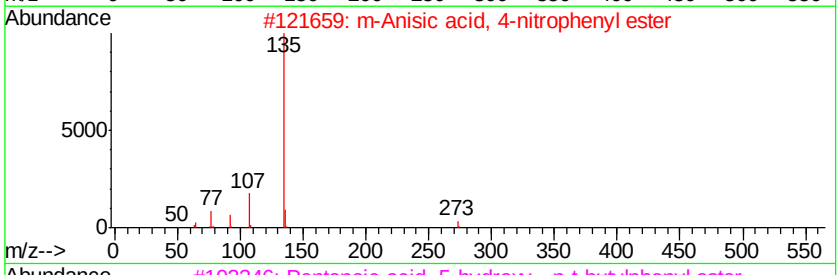
Peak Number 10 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 6

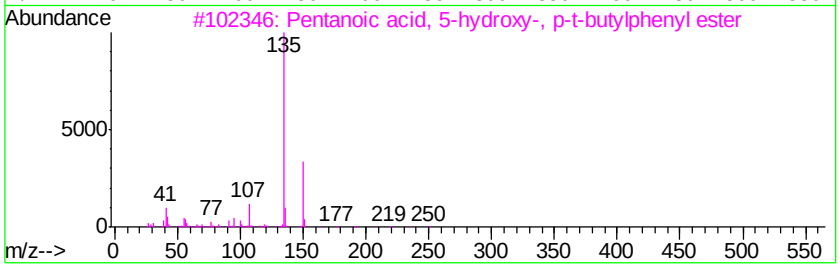
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.03	9.52 ng	1359250	Phenanthrene-d10	11.32

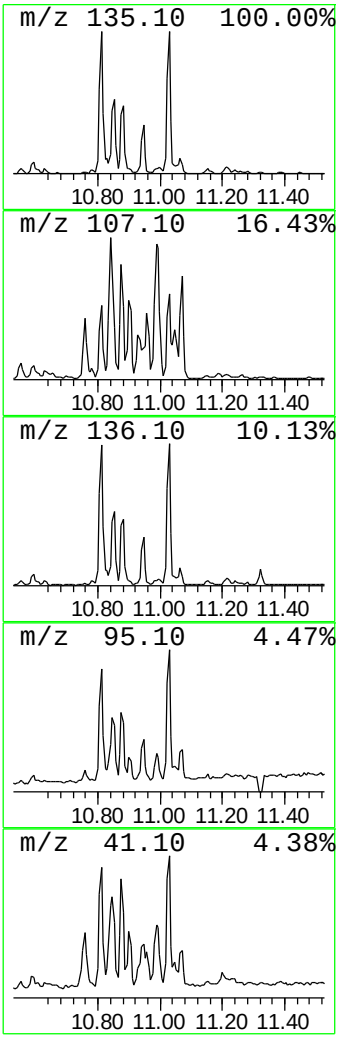
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2	m-Anisic acid, 4-nitrophenyl ester	273	C14H11NO5	1000307-80-4	64
3	Pentanoic acid, 5-hydroxy-, p-t-...	250	C15H22O3	166273-37-6	64
4	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
5	1,2-Benzenediol, o-(4-methoxyben...	362	C22H18O5	1000325-99-0	59











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Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.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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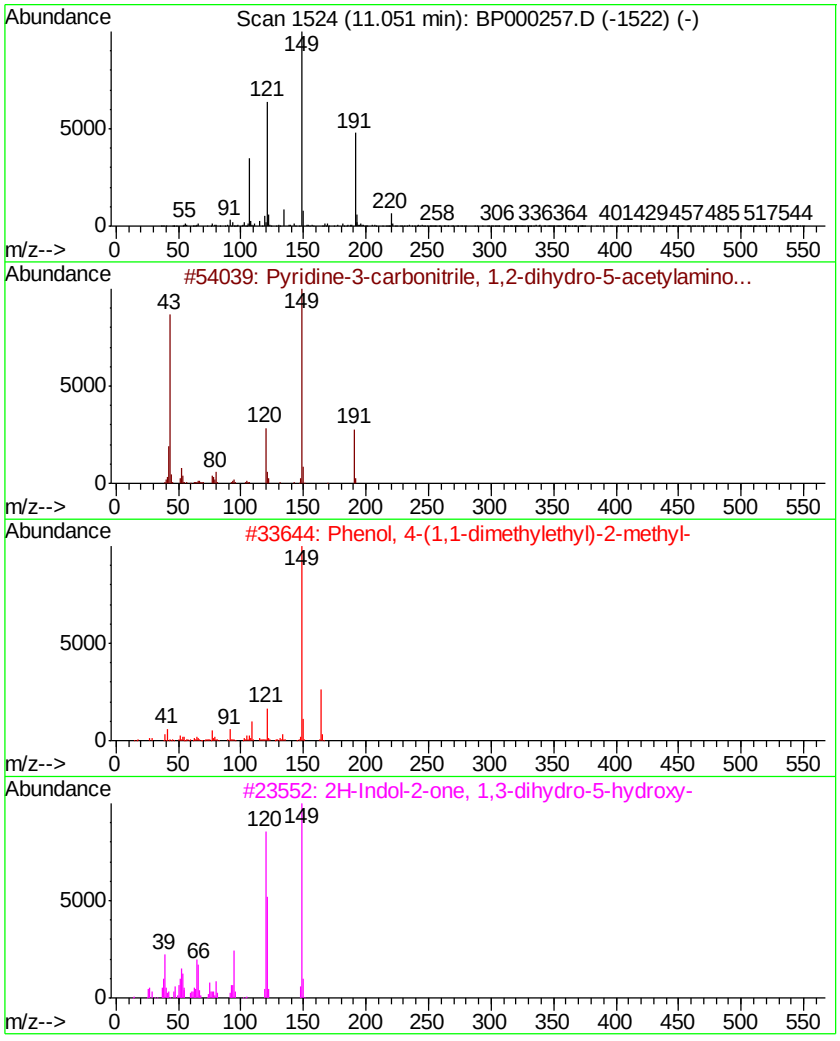
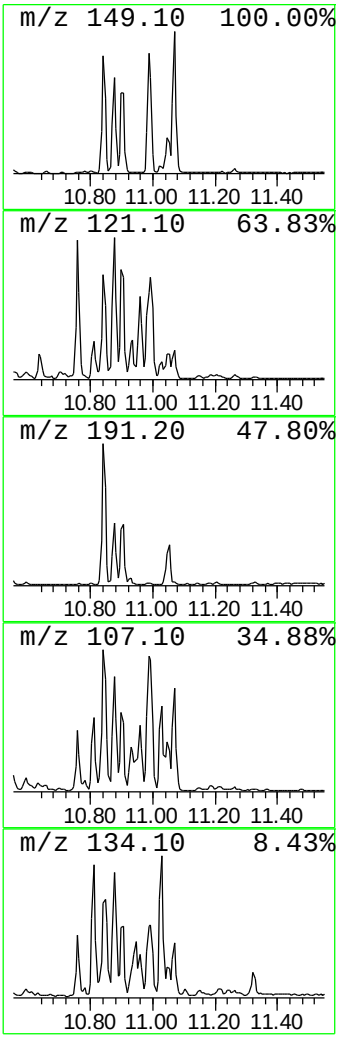
Peak Number 11 Pyridine-3-carbonitrile, 1,... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	2.52 ng	360094	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	52
2	Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	47
3	2H-Indol-2-one, 1,3-dihydro-5-hy...	149	C8H7NO2	003416-18-0	47
4	Phthalic acid, 2-methylallyl pen...	290	C17H22O4	1000315-82-4	43
5	1,2-Benzenedicarboxylic acid, bi...	250	C14H18O4	000605-45-8	43





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.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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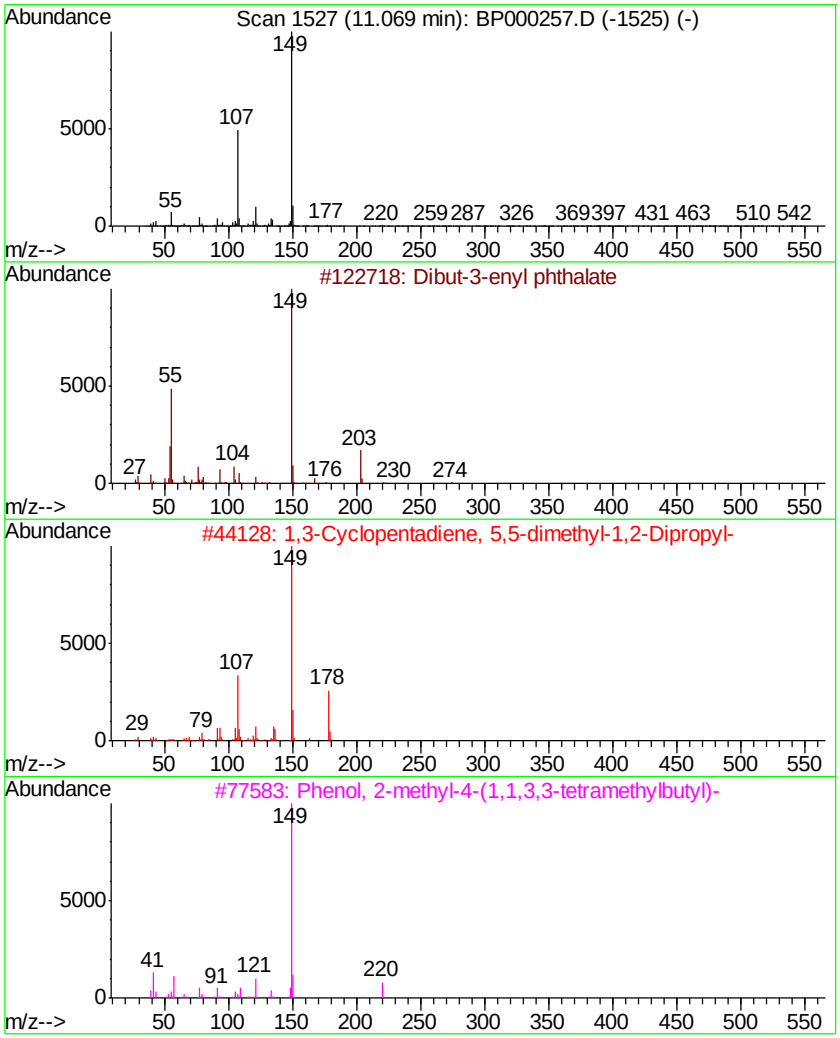
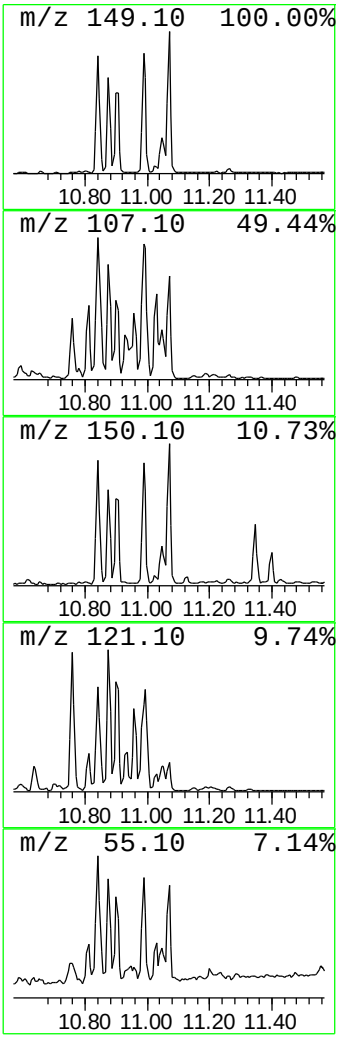
Peak Number 12 unknown11.07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.07	5.02 ng	716634	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibut-3-enyl phthalate	274	C16H18O4	1000373-49-6	47
2	1,3-Cyclopentadiene, 5,5-dimethy...	178	C13H22	1000163-88-0	45
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	40
4	Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	40
5	Phthalic acid, butyl 2-propylphe...	340	C21H24O4	1000357-01-9	40





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
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Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

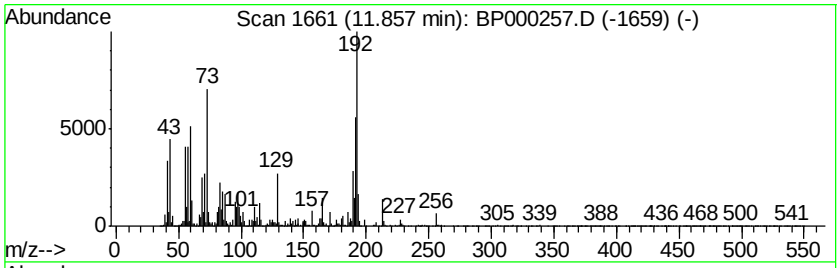
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

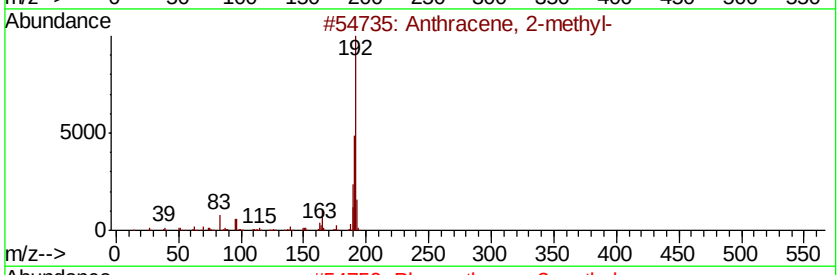
****

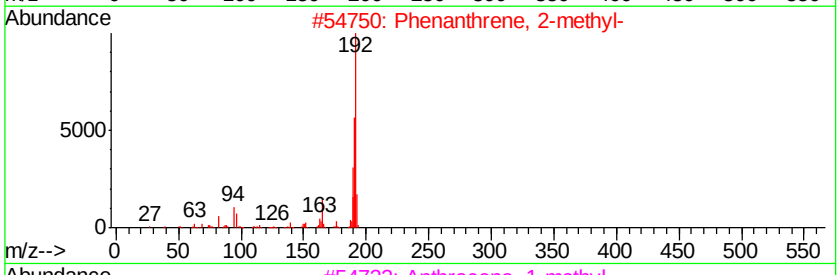
Peak Number 13 Anthracene, 2-methyl- Concentration Rank 13

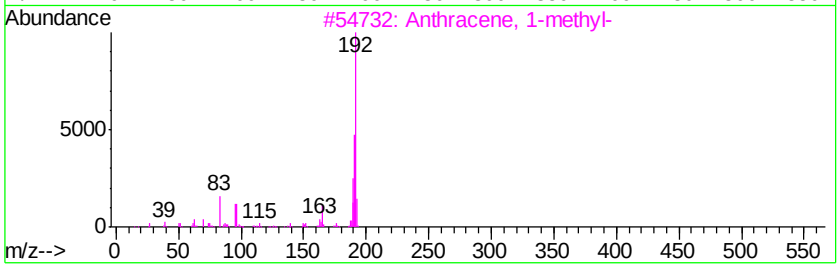
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	3.02 ng	430901	Phenanthrene-d10	11.32

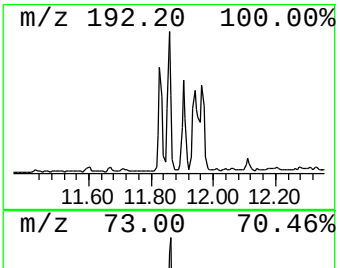
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	92
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92

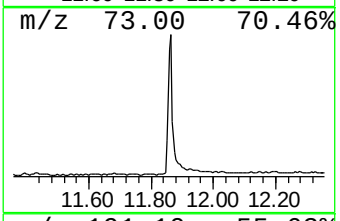


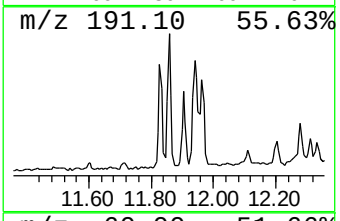


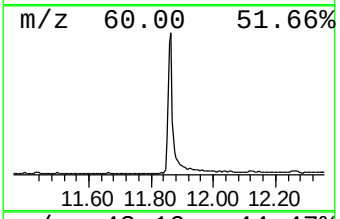


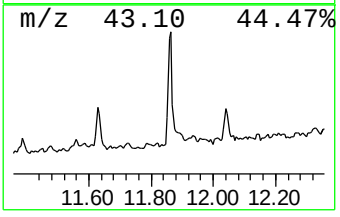












Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-29

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

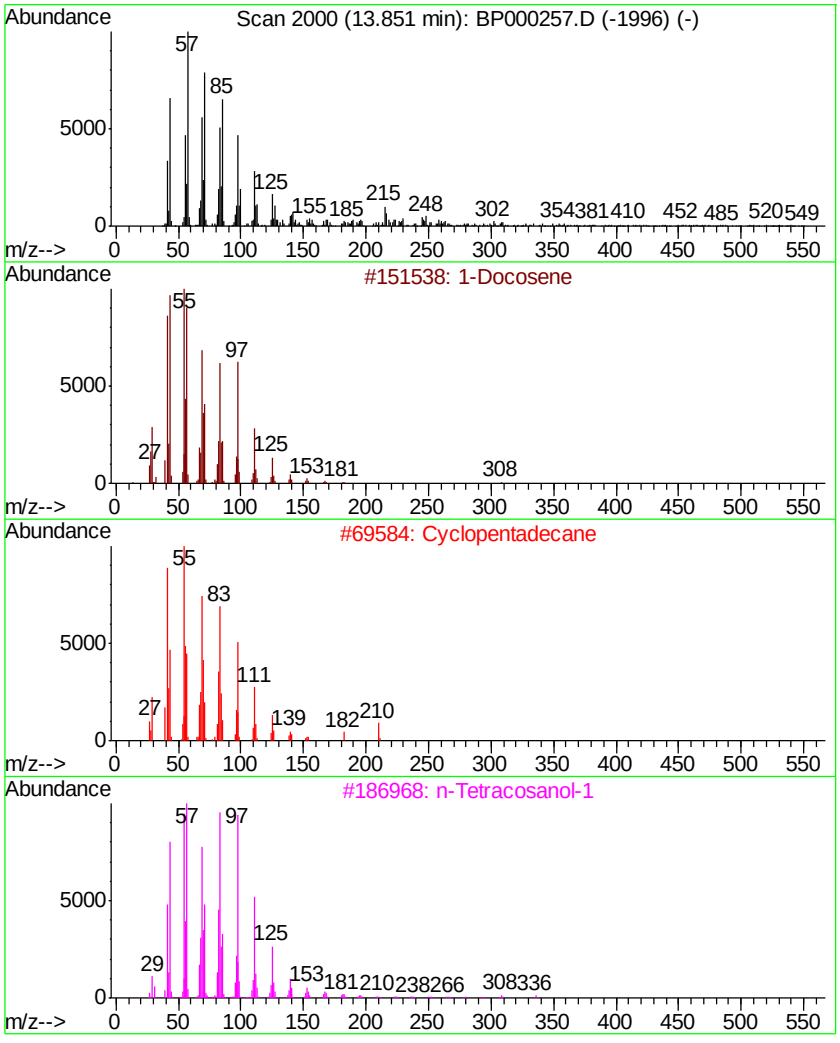
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

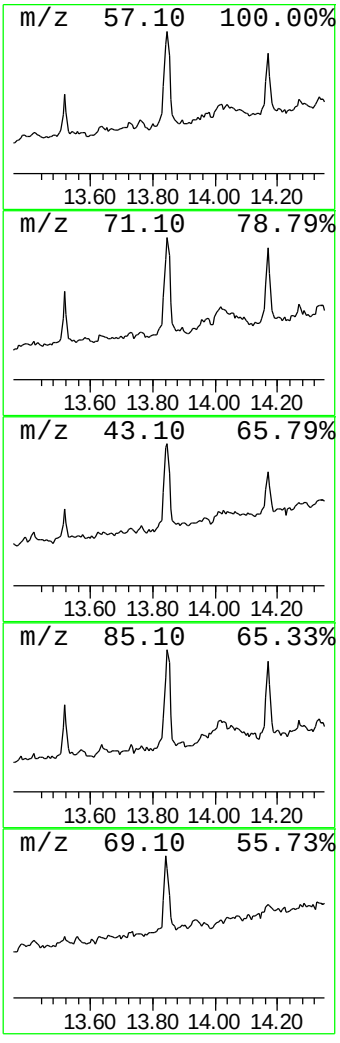
****

Peak Number 14 1-Docosene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	3.27 ng	565495	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	Cyclopentadecane	210	C15H30	000295-48-7	93
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5	17-Pentatriacontene	491	C35H70	006971-40-0	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091619\
Data File : BP000257.D
Acq On : 16 Sep 2019 23:37
Operator : HP/JU
Sample : K4873-04 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_P
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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
unknown6.56	6.56	34.3	ng	3244140	1	6.79	1889640	20.0
unknown10.76	10.76	4.8	ng	683459	4	11.32	2854570	20.0
((1,2-Diethylethy...	10.85	12.4	ng	1772220	4	11.32	2854570	20.0
Phenol, p-tert-bu...	10.88	10.2	ng	1451580	4	11.32	2854570	20.0
Hexestrol dimethy...	10.90	4.8	ng	692048	4	11.32	2854570	20.0
m-Anisic acid, 4-...	10.95	5.5	ng	779919	4	11.32	2854570	20.0
unknown10.99	10.99	9.7	ng	1383470	4	11.32	2854570	20.0
Phenol, 4-(1,1,3,...	11.03	9.5	ng	1359250	4	11.32	2854570	20.0
Pyridine-3-carbon...	11.05	2.5	ng	360094	4	11.32	2854570	20.0
unknown11.07	11.07	5.0	ng	716634	4	11.32	2854570	20.0
Anthracene, 2-met...	11.86	3.0	ng	430901	4	11.32	2854570	20.0
1-Docosene	13.85	3.3	ng	565495	5	13.99	3458170	20.0