LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

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:\HPCHEM1\BNA_F\METHODS\8270-BF010915.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	1.784	57	61	65	rBV	366823	624285	2.48%	0.739%
2	2.310	99	107	110	rVB	8034330	25217068	100.00%	29.864%
3	4.322	278	283	287	rVB	1129832	2315729	9.18%	2.742%
4	4.722	312	318	321	rBV	2244510	5303365	21.03%	6.281%
5	4.984	333	341	346	rVB	3341786	10384653	41.18%	12.298%
6	5.144	352	355	360	rBV	276049	559663	2.22%	0.663%
7	5.499	381	386	389	rVB	1755181	3412018	13.53%	4.041%
8	6.219	445	449	455	rVB	620463	1091678	4.33%	1.293%
9	6.630	482	485	489	rBV	245644	443635	1.76%	0.525%
10	7.030	516	520	522	rBV	195034	353424	1.40%	0.419%
11	7.202	531	535	537	rBV	164923	267328	1.06%	0.317%
12	7.362	545	549	553	rVB2	211276	528069	2.09%	0.625%
13	7.476	553	559	563	rBV3	182313	704632	2.79%	0.834%
14	7.556	563	566	572	rVV	590314	1229283	4.87%	1.456%
15	7.693	572	578	581	rVB2	680480	1593450	6.32%	1.887%
16	7.899	586	596	599	rBV2	172343	425038	1.69%	0.503%
17	7.968	599	602	605	rBV	165708	265959	1.05%	0.315%
18	8.196	619	622	625	rVB	278693	434654	1.72%	0.515%
19	8.379	635	638	641	rVV	572519	1082521	4.29%	1.282%
20	8.448	642	644	646	rVV	243973	374619	1.49%	0.444%
21	8.505	646	649	652	rVB	209527	342080	1.36%	0.405%
22	8.619	652	659	662	rBV	797884	1472952	5.84%	1.744%
23	8.893	676	683	688	rVB	978981	2013368	7.98%	2.384%
24	9.213	708	711	716	rBV3	134239	343409	1.36%	0.407%
25	9.351	719	723	726	rVV	533573	995449	3.95%	1.179%
26	9.419	726	729	734	rVB3	112556	313841	1.24%	0.372%
27	9.705	751	754	757	rBV	314721	496620	1.97%	0.588%
28	10.071	783	786	789	rVB	138988	283248	1.12%	0.335%
29	10.425	812	817	818	rBV2	247941	531935	2.11%	0.630%
30	10.459	818	820	826	rVB2	206672	446630	1.77%	0.529%
31	10.699	836	841	843	rBV3	110891	263256	1.04%	0.312%
32	10.768	843	847	850	rVV	312428	677660	2.69%	0.803%
33	10.905	854	859	861	rVV4	97011	329963	1.31%	0.391%
34	10.962	861	864	866	rVV2	246471	450382	1.79%	0.533%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

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:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.008	866	868	872	rVB	473498	803681	3.19%	0.952%
36	11.202	880	885	888	rVV2	277354	670801	2.66%	0.794%
37	11.499	906	911	915	rBV	228524	459349	1.82%	0.544%
38	12.288	976	980	983	rVV	455470	827692	3.28%	0.980%
39	12.437	987	993	994	rVV2	121893	374784	1.49%	0.444%
40	12.482	994	997	1000	rVV	1009345	1757058	6.97%	2.081%
41	12.574	1001	1005	1008	rVB	588765	987236	3.91%	1.169%
42	12.665	1008	1013	1019	rBV	130128	324572	1.29%	0.384%
43	13.088	1044	1050	1053	rVB	369709	636143	2.52%	0.753%
44	13.191	1053	1059	1062	rBV	122635	303494	1.20%	0.359%
45	13.602	1090	1095	1099	rVB	871207	1603202	6.36%	1.899%
46	13.762	1106	1109	1111	rBV	141680	268432	1.06%	0.318%
47	14.905	1200	1209	1211	rBV	297242	1164641	4.62%	1.379%
48	15.088	1221	1225	1229	rVB	244281	440680	1.75%	0.522%
49	16.403	1337	1340	1342	rBV	192489	264021	1.05%	0.313%
50	16.608	1355	1358	1360	rBV	205153	271150	1.08%	0.321%
51	16.734	1367	1369	1372	rVV	691082	1035849	4.11%	1.227%
52	17.054	1393	1397	1399	rVV	217827	398392	1.58%	0.472%
53	17.351	1420	1423	1426	rBV3	291141	444661	1.76%	0.527%
54	17.820	1462	1464	1467	rVB	258352	328288	1.30%	0.389%
55	17.934	1472	1474	1477	rVB	352035	366236	1.45%	0.434%
56	18.552	1526	1528	1531	rVB	165749	255443	1.01%	0.303%
57	18.780	1546	1548	1552	rBV	335857	521272	2.07%	0.617%
58	18.849	1552	1554	1556	rVV	288425	393816	1.56%	0.466%
59	18.894	1556	1558	1560	rVB	625064	687731	2.73%	0.814%
60	19.946	1647	1650	1653	rBV	429884	492397	1.95%	0.583%
61	20.037	1655	1658	1661	rVV	1053096	1627753	6.45%	1.928%
62	21.306	1767	1769	1777	rVB	344491	467767	1.85%	0.554%
63	21.500	1782	1786	1792	rBV	232864	359918	1.43%	0.426%
64	23.021	1916	1919	1922	rVB	238419	332313	1.32%	0.394%

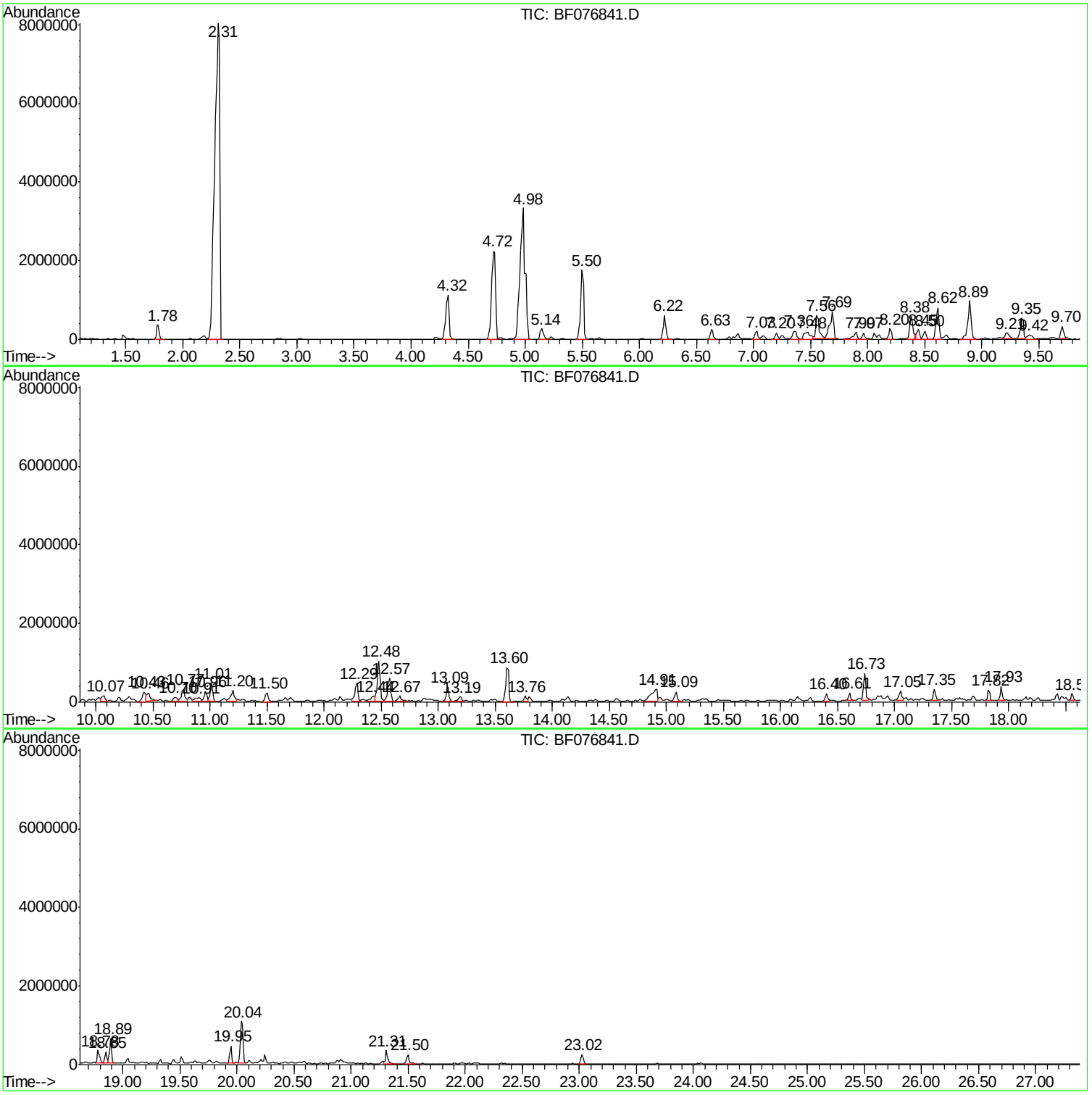
Sum of corrected areas: 84440636

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

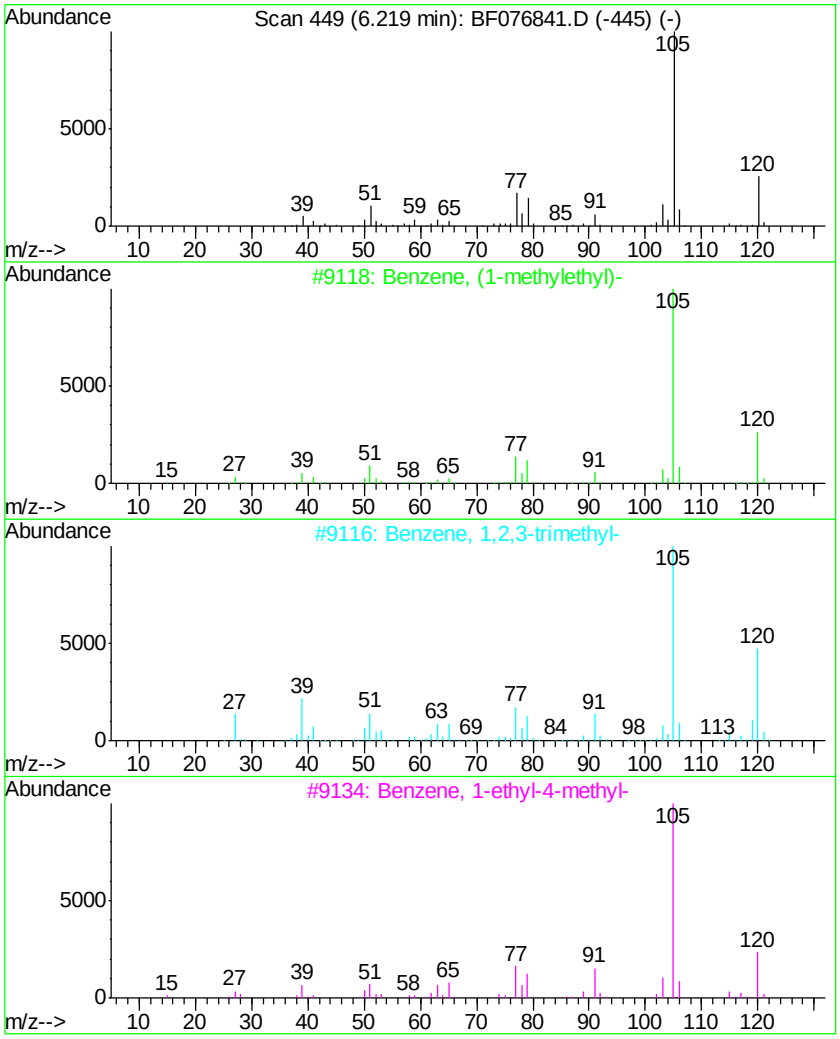
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

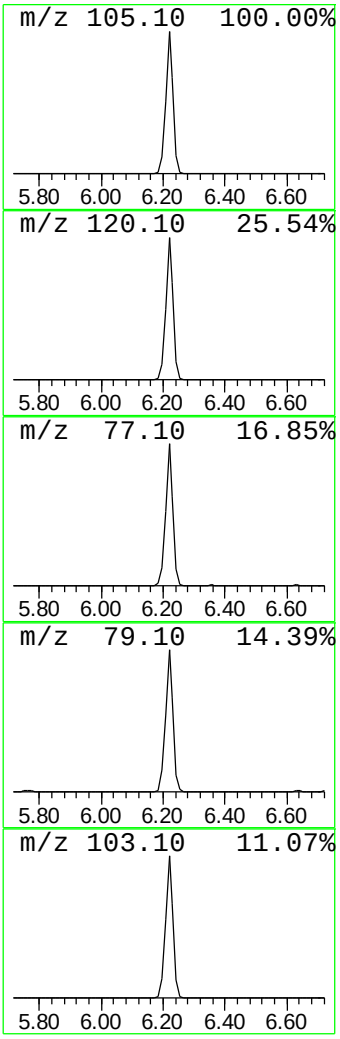
****

Peak Number 7 Benzene, (1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.22	82.09 ng	1091680	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72
4	2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	72
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

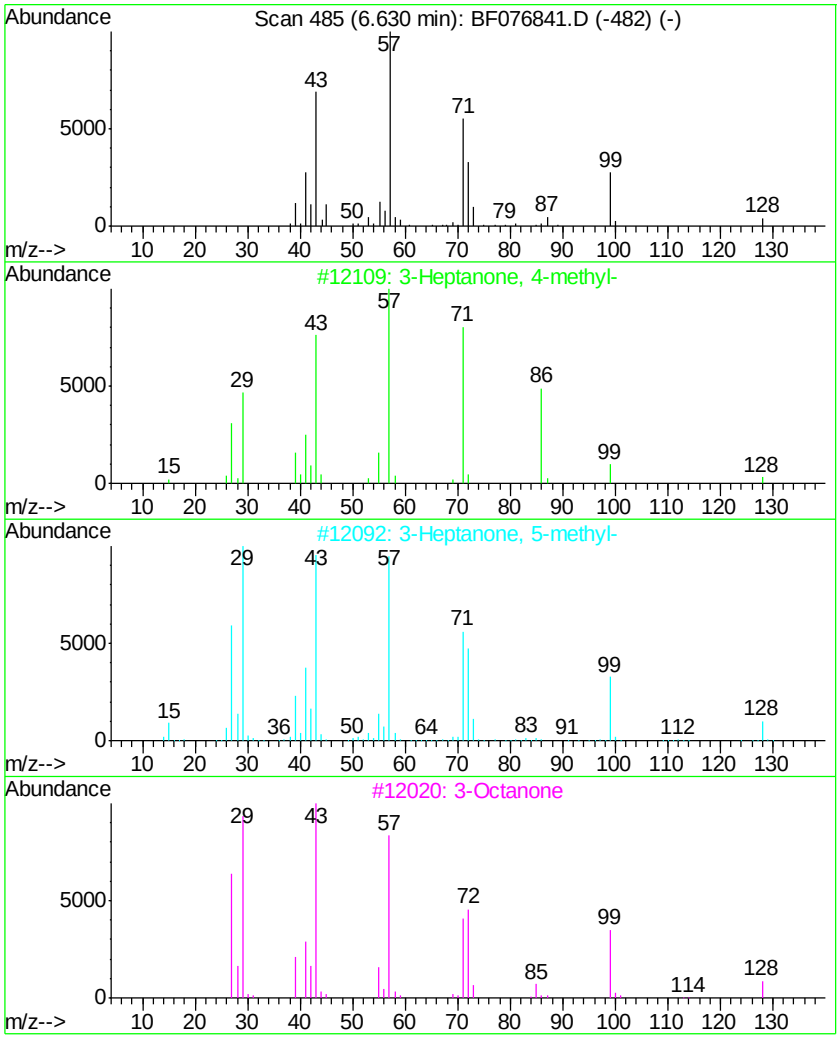
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

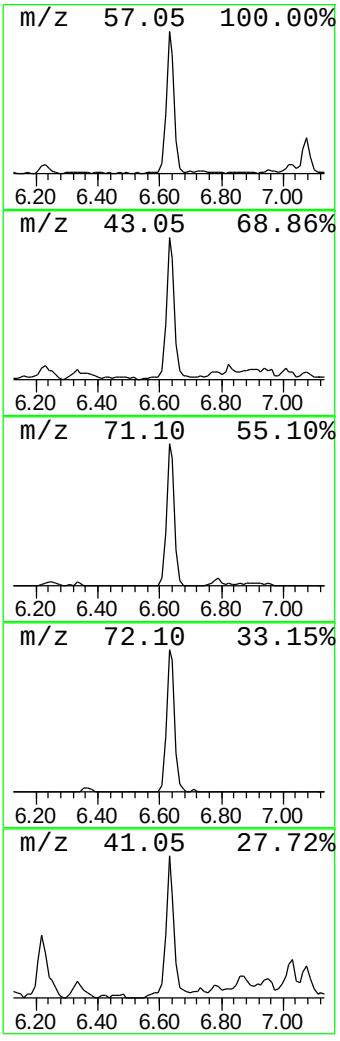
****

Peak Number 8 unknown6.63 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	33.36 ng	443635	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanone, 4-methyl-	128	C8H16O	006137-11-7	46
2	3-Heptanone, 5-methyl-	128	C8H16O	000541-85-5	45
3	3-Octanone	128	C8H16O	000106-68-3	42
4	Hexanoic acid, 1,1-dimethylpropy...	186	C11H22O2	116423-69-9	38
5	Allyl 2-ethyl butyrate	156	C9H16O2	007493-69-8	37





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

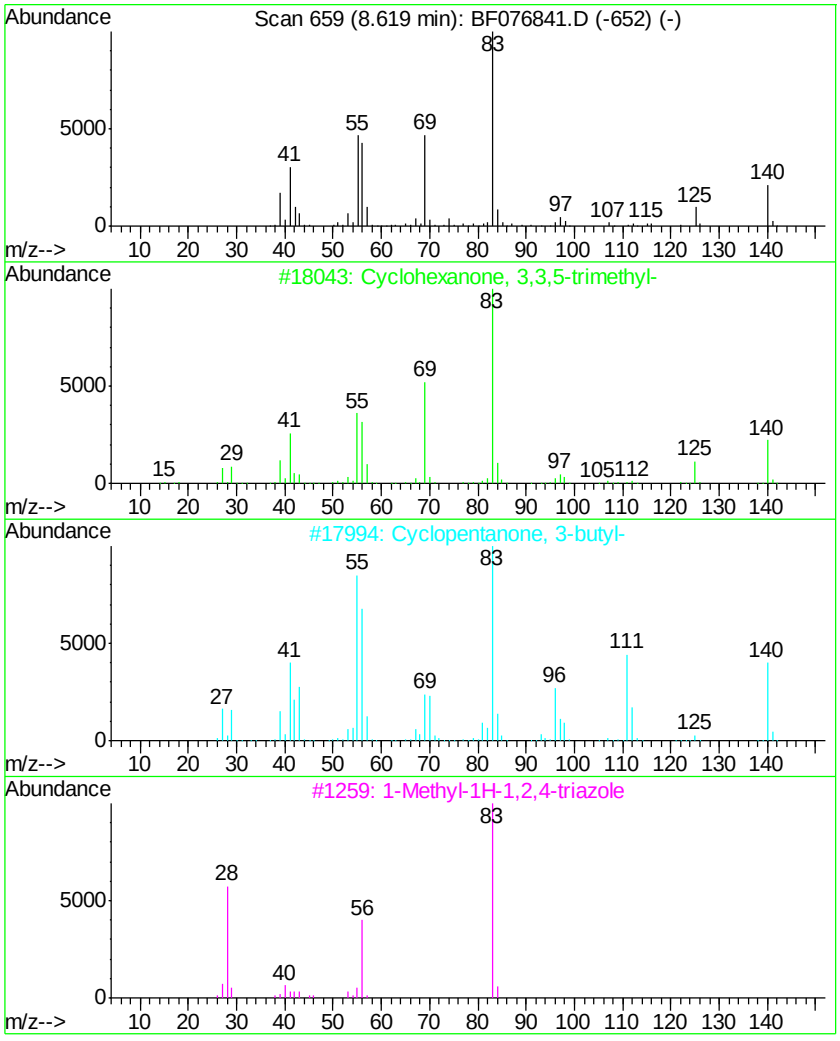
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

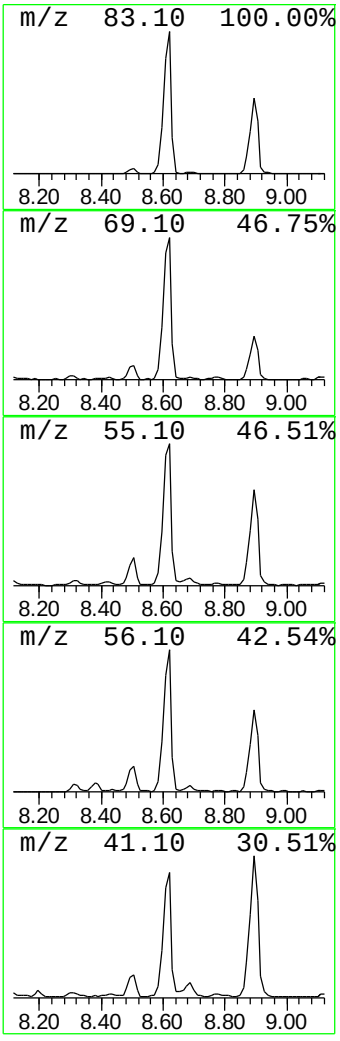
****

Peak Number 11 Cyclohexanone, 3,3,5-trimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	110.77 ng	1472950	1,4-Dichlorobenzene-d4	8.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
2	Cyclopentanone, 3-butyl-	140	C9H16O	057283-81-5	58
3	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	50
4	Cyclohexane, butyl-	140	C10H20	001678-93-9	50
5	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	50





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

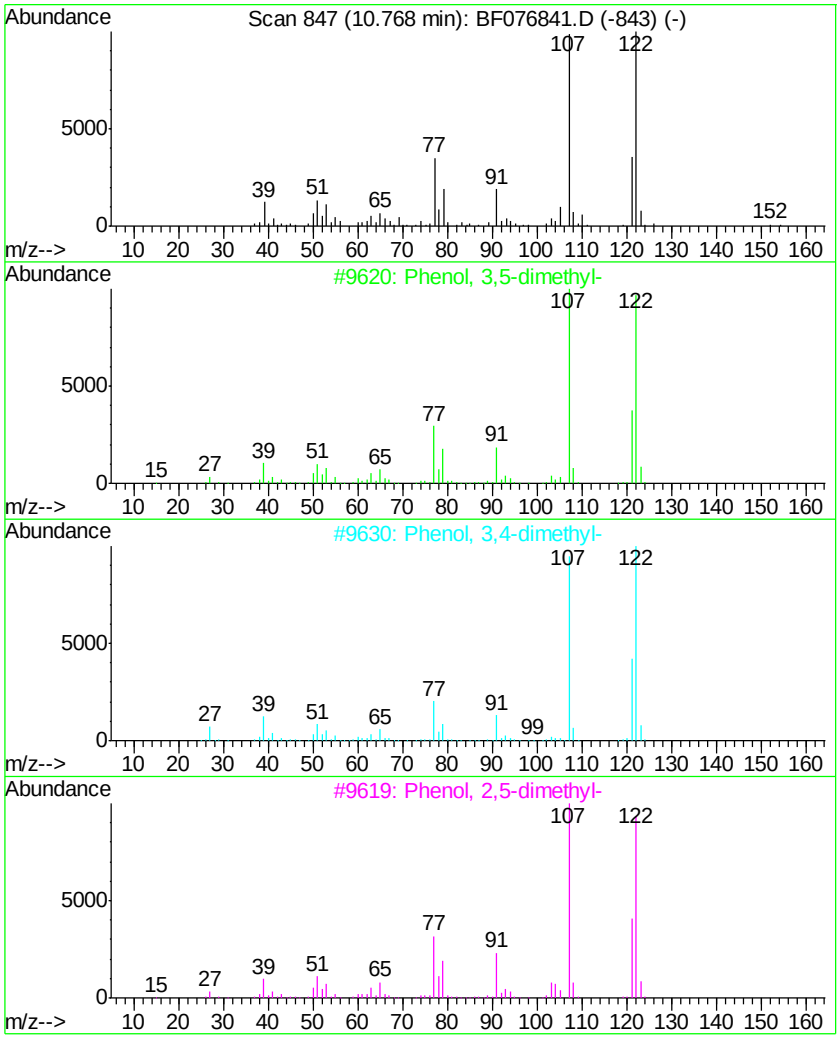
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

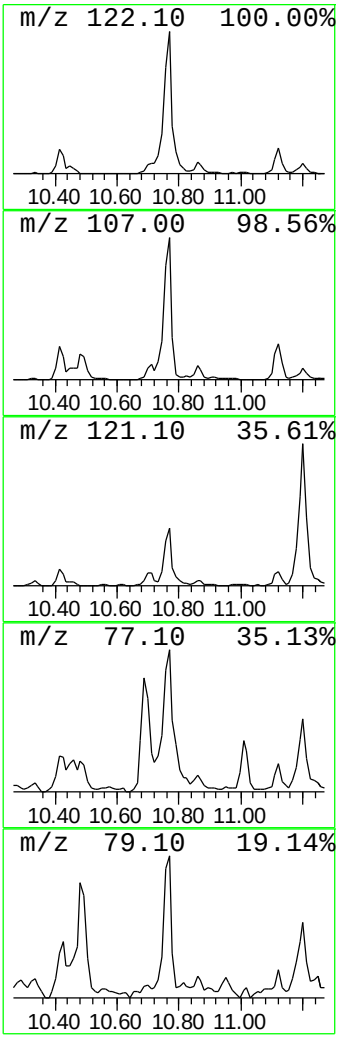
****

Peak Number 12 Phenol, 3,5-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	30.09 ng	677660	Naphthalene-d8	10.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	93
3	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	93
4	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	90
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	90





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

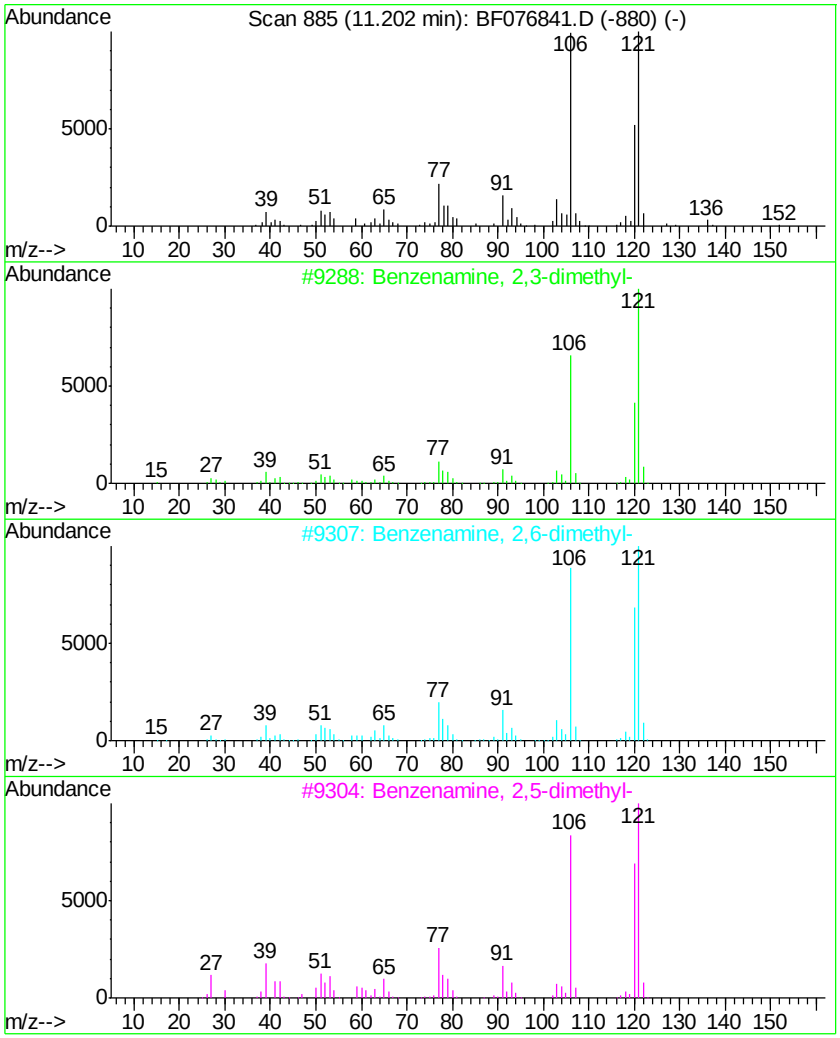
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

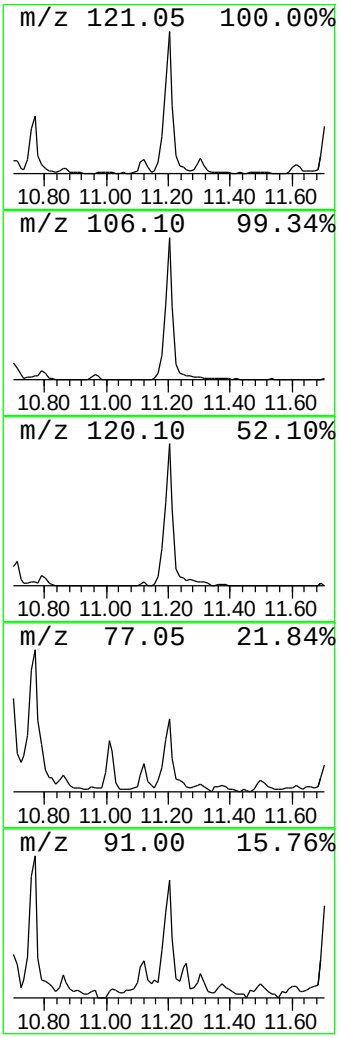
****

Peak Number 13 Benzenamine, 2,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	29.79 ng	670801	Naphthalene-d8	10.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,3-dimethyl-	121	C8H11N	000087-59-2	97
2	Benzenamine, 2,6-dimethyl-	121	C8H11N	000087-62-7	97
3	Benzenamine, 2,5-dimethyl-	121	C8H11N	000095-78-3	96
4	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	96
5	Benzenamine, 3,5-dimethyl-	121	C8H11N	000108-69-0	96





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

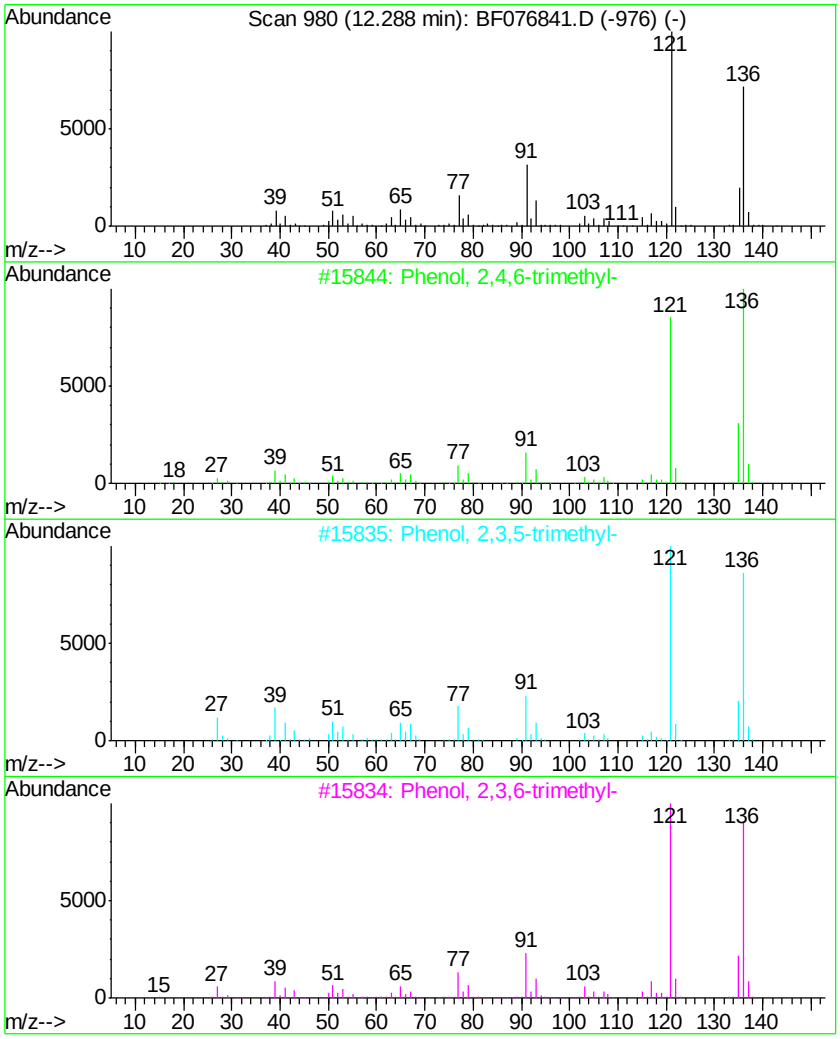
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

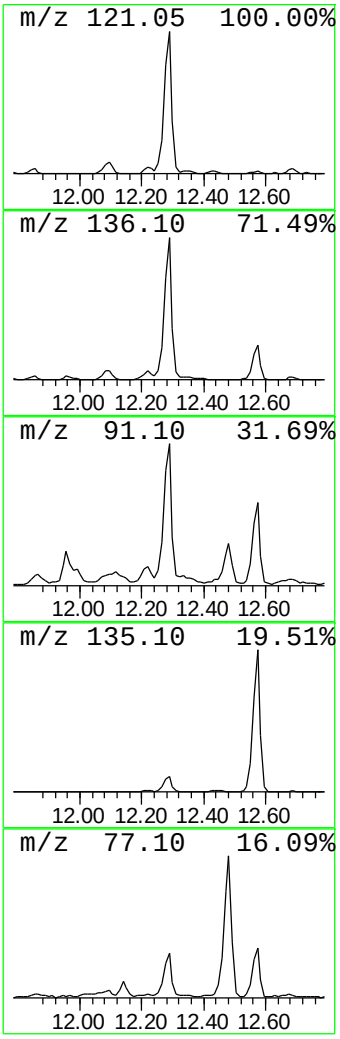
****

Peak Number 14 Phenol, 2,4,6-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	36.76 ng	827692	Naphthalene-d8	10.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	97
2	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	97
3	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	95
4	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	91
5	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	91





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

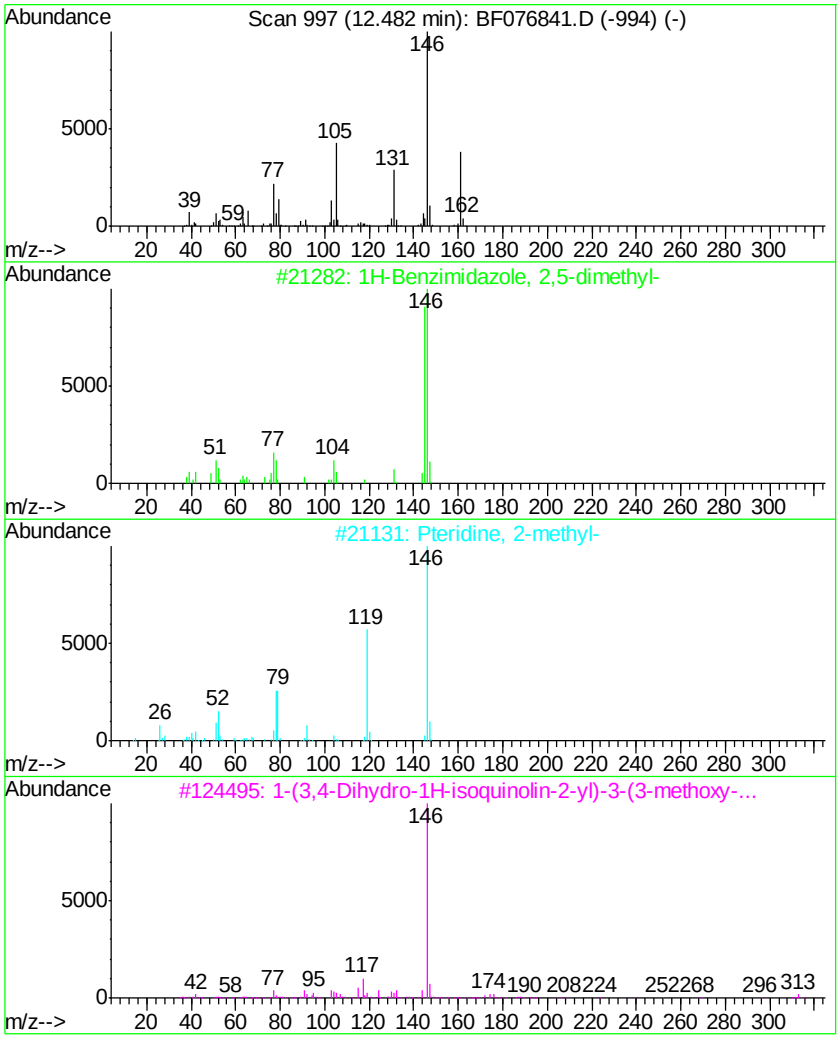
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

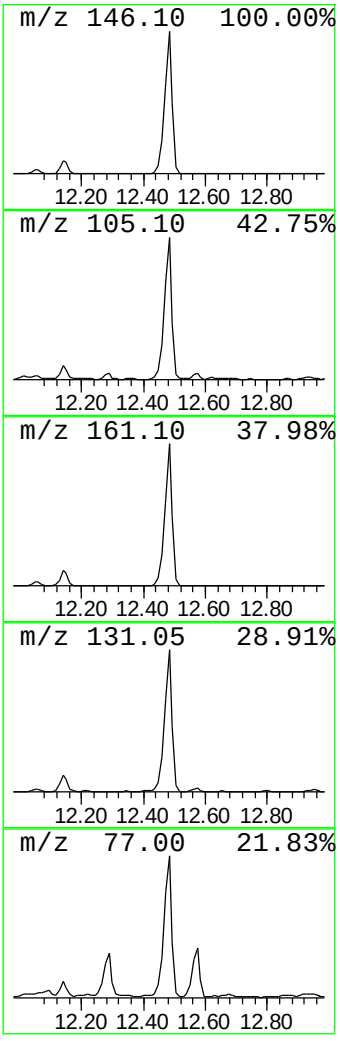
****

Peak Number 15 unknown12.48 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.48	78.03 ng	1757060	Naphthalene-d8	10.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2,5-dimethyl-	146	C9H10N2	001792-41-2	40
2	Pteridine, 2-methyl-	146	C7H6N4	002432-20-4	37
3	1-(3,4-Dihydro-1H-isoquinolin-2-...	313	C19H23NO3	1000275-12-7	37
4	Isoquinoline, 1,2,3,4-tetrahydro...	161	C11H15N	054365-72-9	36
5	1,2,5-Triamino-1H-pyrrole-3,4-di...	162	C6H6N6	075610-90-1	33





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

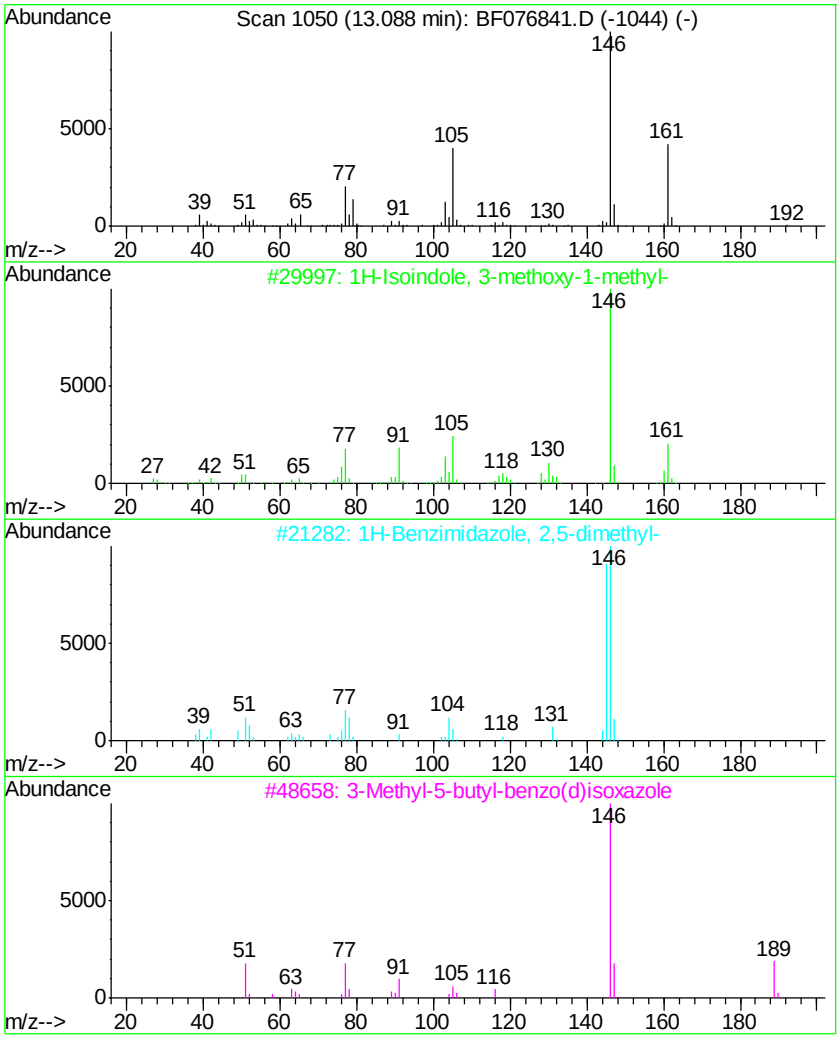
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

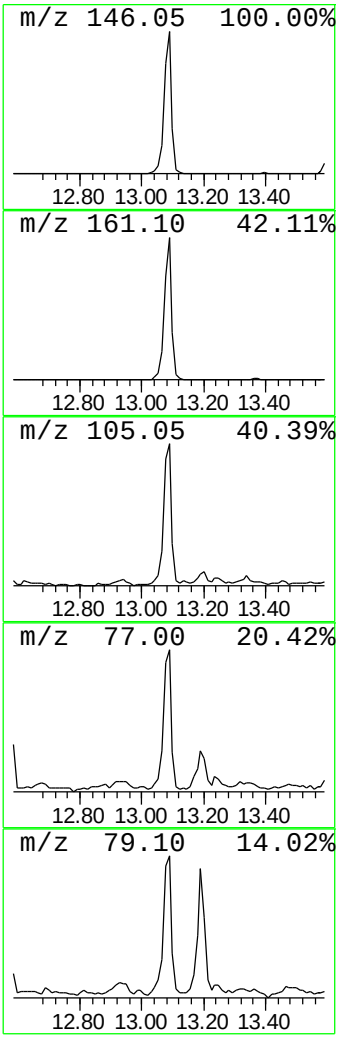
****

Peak Number 17 1H-Isoindole, 3-methoxy-1-m... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.09	28.87 ng	636143	Acenaphthene-d10	15.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Isoindole, 3-methoxy-1-methyl-	161	C10H11NO	083889-38-7	50
2	1H-Benzimidazole, 2,5-dimethyl-	146	C9H10N2	001792-41-2	28
3	3-Methyl-5-butyl-benzo(d)isoxazole	189	C12H15NO	131222-25-8	25
4	2-Methyl-5-butyl-benzo(d)oxazole	189	C12H15NO	131222-26-9	25
5	5,6-Dimethyl-1,3.4.4-dithiapyrid...	146	C4H6N2S2	021423-60-9	25





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

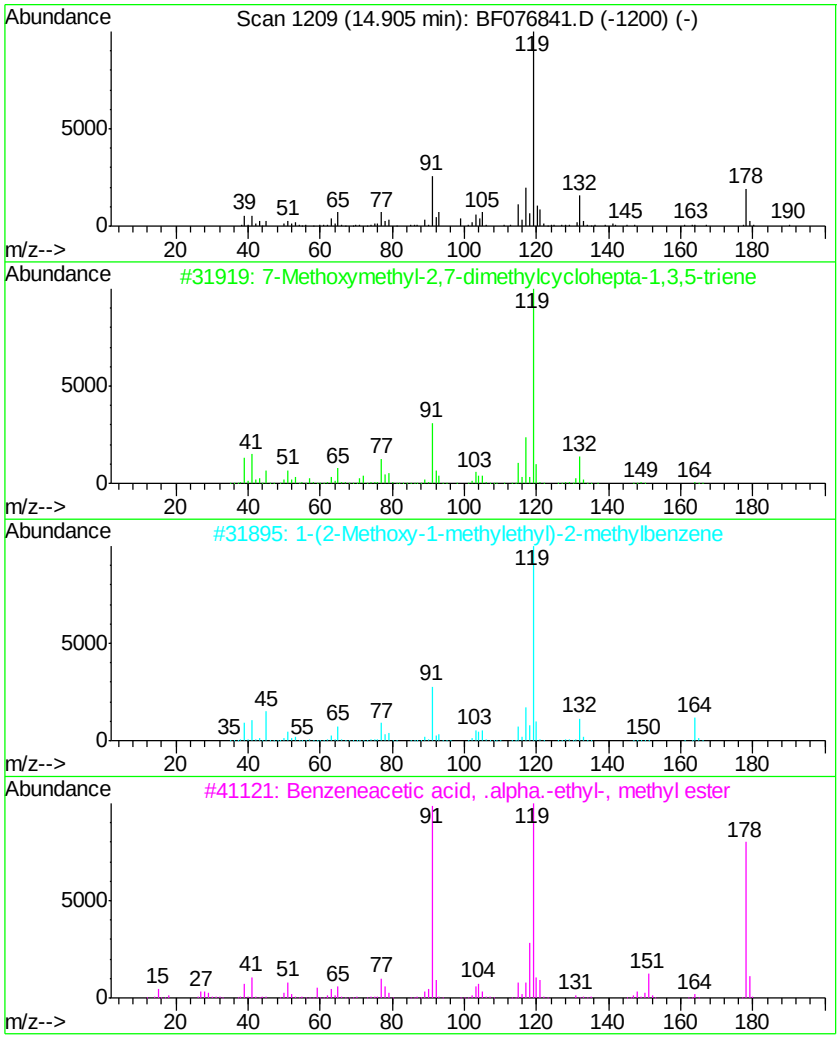
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

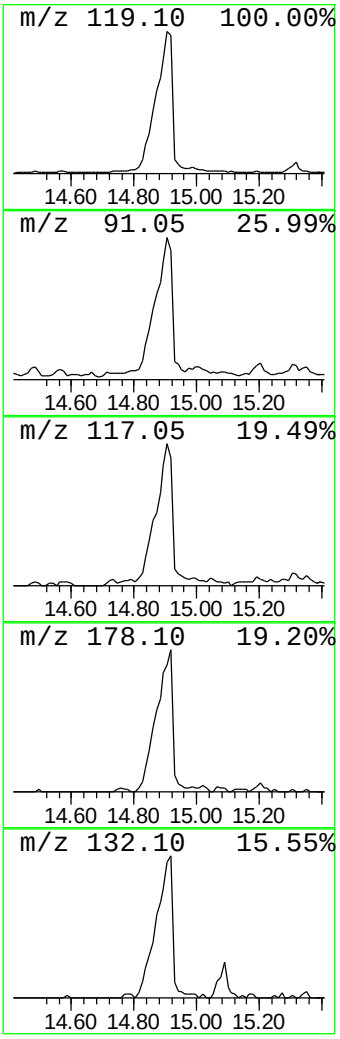
****

Peak Number 18 7-Methoxymethyl-2,7-dimethy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	52.86 ng	1164640	Acenaphthene-d10	15.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methoxymethyl-2,7-dimethylcycl...	164	C11H16O	073992-48-0	81
2	1-(2-Methoxy-1-methylethyl)-2-me...	164	C11H16O	1000210-02-1	72
3	Benzeneacetic acid, .alpha.-ethy...	178	C11H14O2	002294-71-5	64
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	58
5	Benzene, 4-(chloromethyl)-1,2-di...	154	C9H11Cl	000102-46-5	53





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

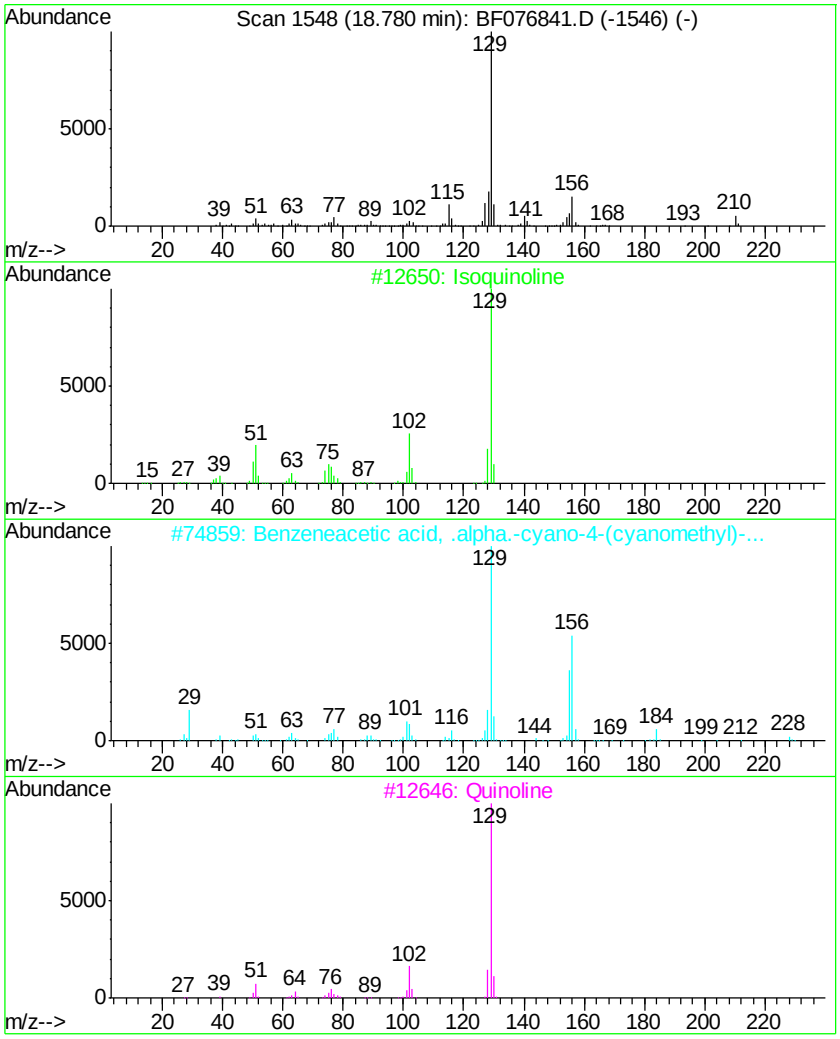
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

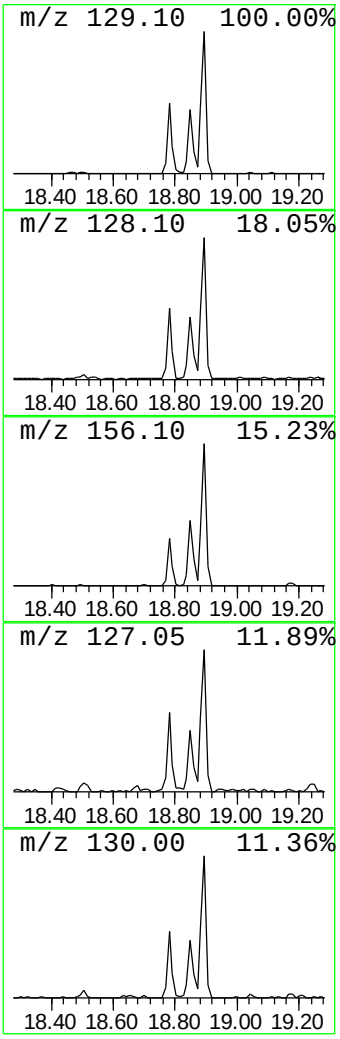
****

Peak Number 19 Isoquinoline Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.78	31.76 ng	521272	Phenanthrene-d10	17.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isoquinoline	129	C9H7N	000119-65-3	53
2	Benzeneacetic acid, .alpha.-cyan...	228	C13H12N2O2	134882-04-5	50
3	Quinoline	129	C9H7N	000091-22-5	40
4	Benzeneacetonitrile, .alpha.-met...	129	C9H7N	000495-10-3	36
5	1,4-Methanonaphthalen-9-ol, 1,2,...	202	C13H14O2	001207-28-9	36





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

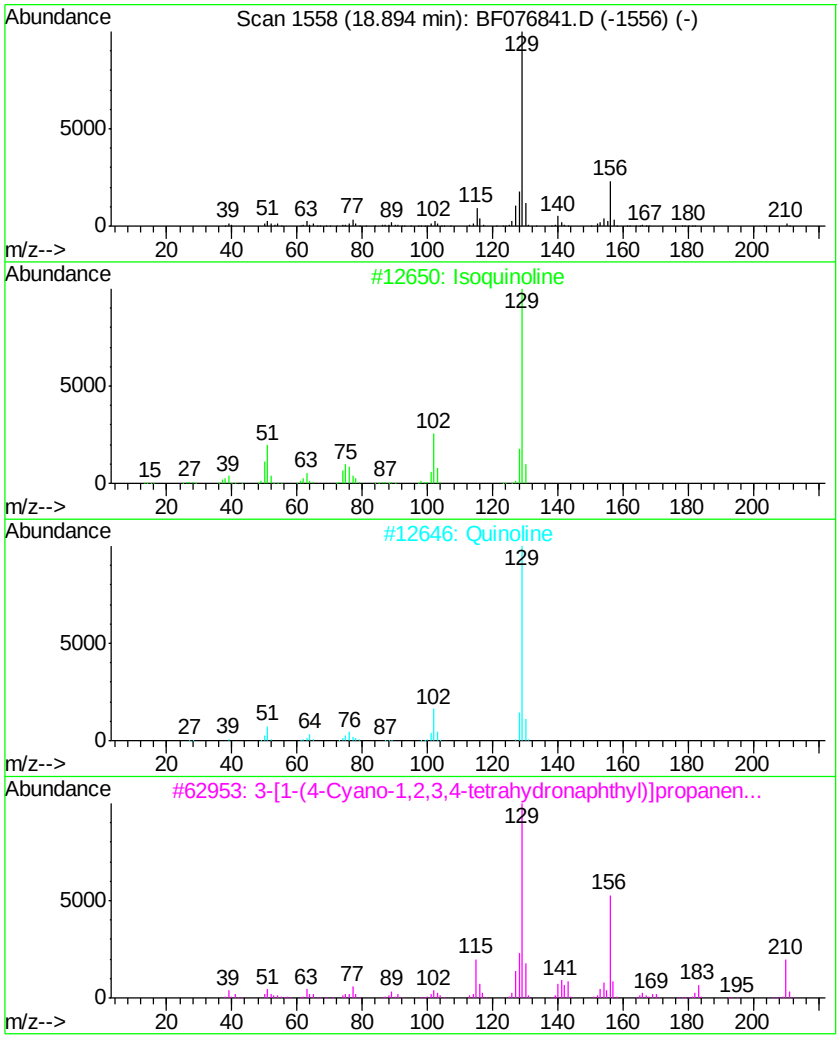
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

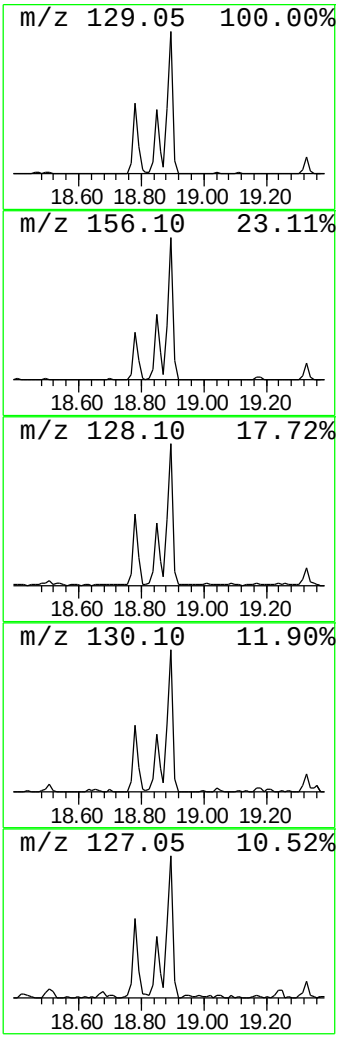
****

Peak Number 20 unknown18.89 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.89	41.90 ng	687731	Phenanthrene-d10	17.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isoquinoline	129	C9H7N	000119-65-3	50
2	Quinoline	129	C9H7N	000091-22-5	49
3	3-[1-(4-Cyano-1,2,3,4-tetrahydro...	210	C14H14N2	057964-40-6	47
4	Benzeneacetonitrile, .alpha.-met...	129	C9H7N	000495-10-3	40
5	2-Propenenitrile, 3-phenyl-, (E)-	129	C9H7N	001885-38-7	40





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011315\
Data File : BF076841.D
Acq On : 13 Jan 2015 13:24
Operator : IZ / TP
Sample : G1044-01
Misc : W
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-1A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, (1-methy...	6.22	82.1	ng	1091680	1	8.06	265959	20.0
unknown6.63	6.63	33.4	ng	443635	1	8.06	265959	20.0
Cyclohexanone, 3,...	8.62	110.8	ng	1472950	1	8.06	265959	20.0
Phenol, 3,5-dimet...	10.77	30.1	ng	677660	2	10.96	450382	20.0
Benzenamine, 2,3-...	11.20	29.8	ng	670801	2	10.96	450382	20.0
Phenol, 2,4,6-tri...	12.29	36.8	ng	827692	2	10.96	450382	20.0
unknown12.48	12.48	78.0	ng	1757060	2	10.96	450382	20.0
1H-Isoindole, 3-m...	13.09	28.9	ng	636143	3	15.09	440680	20.0
7-Methoxymethyl-2...	14.91	52.9	ng	1164640	3	15.09	440680	20.0
Isoquinoline	18.78	31.8	ng	521272	4	17.82	328288	20.0
unknown18.89	18.89	41.9	ng	687731	4	17.82	328288	20.0