LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-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_F\METHODS\8270-BF081821.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125263.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.187	10	17	26	rVB	1348064	1789861	32.84%	3.155%
2	5.534	580	586	589	rBV	4763536	4858880	89.15%	8.564%
3	6.539	751	757	760	rBV	4532798	4851889	89.02%	8.551%
4	6.910	817	820	824	rBV	1182877	1011737	18.56%	1.783%
5	7.475	911	916	919	rBV	3406853	3205820	58.82%	5.650%
6	8.092	1018	1021	1032	rBV	1044008	1268199	23.27%	2.235%
7	8.169	1032	1034	1036	rVV	100705	79306	1.46%	0.140%
8	8.192	1036	1038	1041	rVV	1477595	1225515	22.49%	2.160%
9	8.486	1085	1088	1093	rBV2	42360	58844	1.08%	0.104%
10	8.610	1105	1109	1111	rBV	118899	97524	1.79%	0.172%
11	8.780	1134	1138	1141	rBV	390799	313104	5.74%	0.552%
12	9.269	1217	1221	1224	rBV	6000733	5373185	98.59%	9.470%
13	9.345	1230	1234	1237	rBV	270264	224176	4.11%	0.395%
14	9.380	1237	1240	1244	rVB3	71581	72459	1.33%	0.128%
15	9.875	1320	1324	1328	rVB	71426	64537	1.18%	0.114%
16	9.951	1333	1337	1340	rVV	1817475	1524663	27.97%	2.687%
17	9.980	1340	1342	1346	rVB	174242	130852	2.40%	0.231%
18	10.151	1366	1371	1375	rBV	108982	118090	2.17%	0.208%
19	10.498	1426	1430	1435	rVB	135864	133510	2.45%	0.235%
20	10.739	1467	1471	1474	rBV	4437021	3879245	71.18%	6.837%
21	10.857	1486	1491	1494	rBV2	64636	82634	1.52%	0.146%
22	11.045	1517	1523	1526	rBV4	46734	82315	1.51%	0.145%
23	11.239	1553	1556	1560	rVB	113339	95184	1.75%	0.168%
24	11.292	1560	1565	1567	rVV3	62286	78984	1.45%	0.139%
25	11.333	1567	1572	1577	rVB2	114976	149708	2.75%	0.264%
26	11.439	1586	1590	1592	rBV	1743187	1757244	32.24%	3.097%
27	11.463	1592	1594	1597	rVV	1685081	1323544	24.28%	2.333%
28	11.510	1599	1602	1605	rVB	369070	305433	5.60%	0.538%
29	11.663	1623	1628	1636	rBV2	210364	262204	4.81%	0.462%
30	11.727	1636	1639	1643	rVV4	44904	63566	1.17%	0.112%
31	11.821	1652	1655	1657	rBV	160192	114613	2.10%	0.202%
32	11.957	1671	1678	1683	rBV2	645148	979749	17.98%	1.727%
33	12.010	1685	1687	1690	rVV	87998	88160	1.62%	0.155%
34	12.051	1690	1694	1696	rVV2	323268	341193	6.26%	0.601%
35	12.069	1696	1697	1701	rVB	126008	101705	1.87%	0.179%
36	12.127	1702	1707	1712	rBV4	47886	85289	1.56%	0.150%

LSC Area Percent Report

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Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-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_F\METHODS\8270-BF081821.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.233	1722	1725	1727	rBV	150130	134510	2.47%	0.237%
38	12.257	1727	1729	1733	rVB	176806	141618	2.60%	0.250%
39	12.486	1765	1768	1771	rBV2	108133	125759	2.31%	0.222%
40	12.527	1771	1775	1780	rVV4	157632	224694	4.12%	0.396%
41	12.574	1780	1783	1787	rVB	103105	117651	2.16%	0.207%
42	12.651	1792	1796	1799	rBV	2388471	2010974	36.90%	3.544%
43	12.739	1808	1811	1815	rBV2	290201	278701	5.11%	0.491%
44	12.804	1819	1822	1824	rVB	67648	64248	1.18%	0.113%
45	12.880	1829	1835	1840	rVV	1845791	1994952	36.60%	3.516%
46	12.927	1840	1843	1847	rVB2	96504	116441	2.14%	0.205%
47	13.021	1855	1859	1866	rVB	5203516	5450158	100.00%	9.606%
48	13.098	1867	1872	1876	rBV2	108109	203351	3.73%	0.358%
49	13.180	1883	1886	1888	rBV4	89649	111381	2.04%	0.196%
50	13.204	1888	1890	1894	rVB	237707	213684	3.92%	0.377%
51	13.245	1894	1897	1899	rBV3	75256	85480	1.57%	0.151%
52	13.274	1899	1902	1904	rBV2	168057	141212	2.59%	0.249%
53	13.310	1904	1908	1911	rVV2	160199	159762	2.93%	0.282%
54	13.404	1921	1924	1926	rBV	76876	71815	1.32%	0.127%
55	13.433	1927	1929	1932	rBV2	81280	75555	1.39%	0.133%
56	13.586	1952	1955	1957	rBV2	85920	89563	1.64%	0.158%
57	13.715	1974	1977	1981	rVB	161072	168283	3.09%	0.297%
58	13.827	1992	1996	1998	rBV2	128038	141859	2.60%	0.250%
59	13.857	1998	2001	2003	rVV	119189	115140	2.11%	0.203%
60	13.880	2003	2005	2009	rVV2	107972	153673	2.82%	0.271%
61	13.921	2009	2012	2016	rVB2	196887	192682	3.54%	0.340%
62	14.074	2031	2038	2041	rBV3	1509872	2131901	39.12%	3.757%
63	14.104	2041	2043	2049	rVB	826701	681238	12.50%	1.201%
64	14.180	2054	2056	2059	rBV	96509	83761	1.54%	0.148%
65	14.462	2102	2104	2107	rVB	150860	116057	2.13%	0.205%
66	14.498	2108	2110	2113	rVB2	84469	69574	1.28%	0.123%
67	14.539	2115	2117	2119	rVV	71254	63163	1.16%	0.111%
68	14.592	2123	2126	2127	rBV	79698	73412	1.35%	0.129%
69	14.933	2181	2184	2186	rBV	120070	131481	2.41%	0.232%
70	15.133	2214	2218	2221	rBV	705032	1056164	19.38%	1.861%
71	15.245	2234	2237	2240	rVB	131536	132445	2.43%	0.233%
72	15.445	2268	2271	2277	rVB	413644	568017	10.42%	1.001%
73	15.509	2278	2282	2288	rBV	591311	716138	13.14%	1.262%
74	15.580	2289	2294	2297	rBV	943558	1232199	22.61%	2.172%
75	17.086	2545	2550	2558	rVB2	265508	547129	10.04%	0.964%
76	17.545	2623	2628	2638	rVB	184509	360498	6.61%	0.635%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-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_F\METHODS\8270-BF081821.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

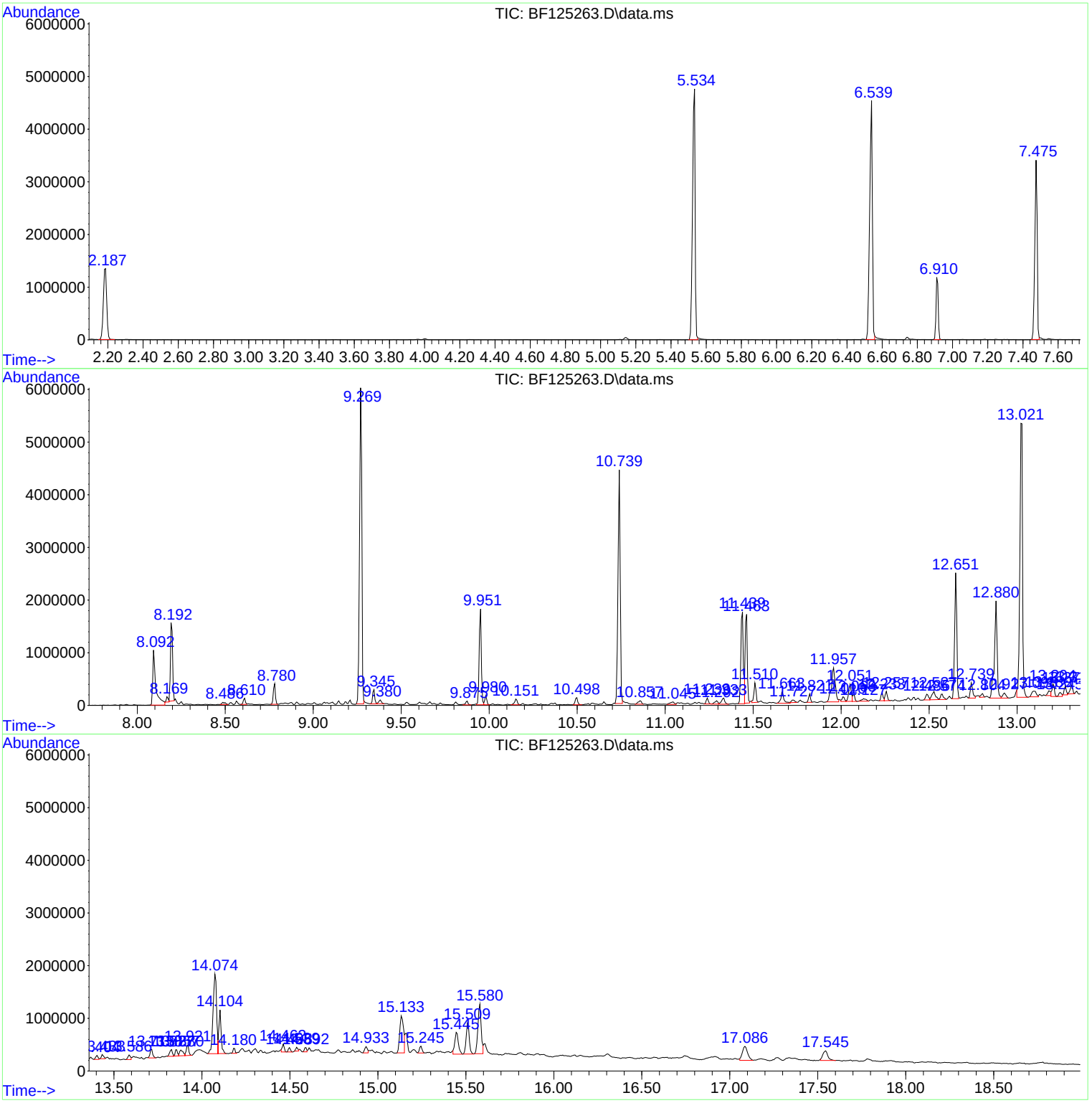
Sum of corrected areas: 56739244

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

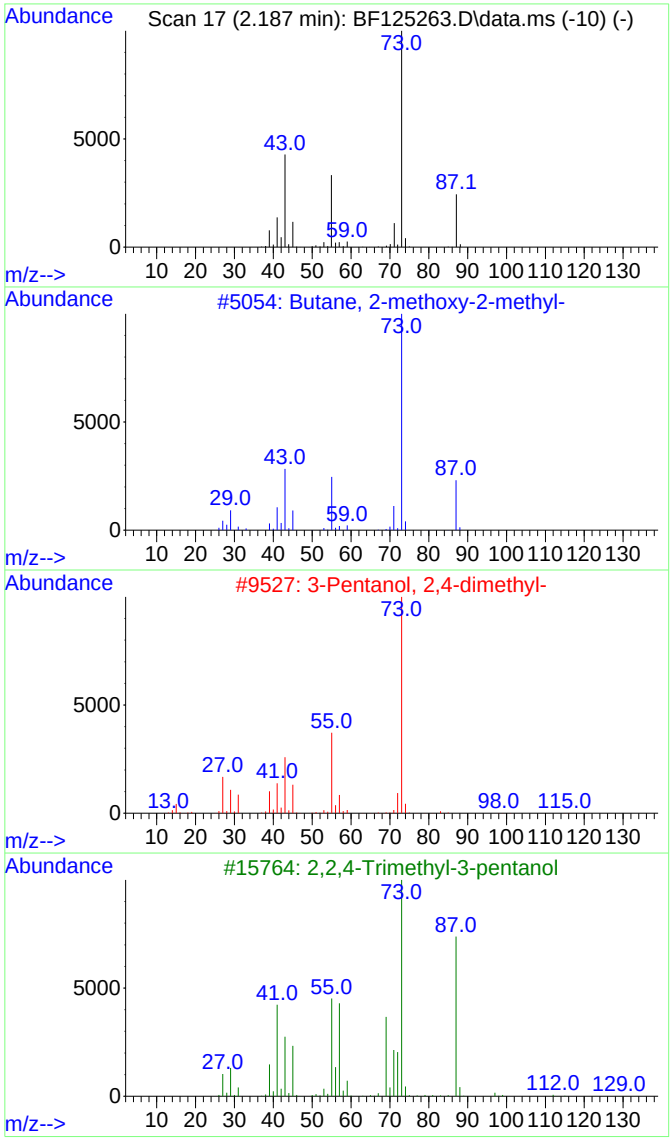
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

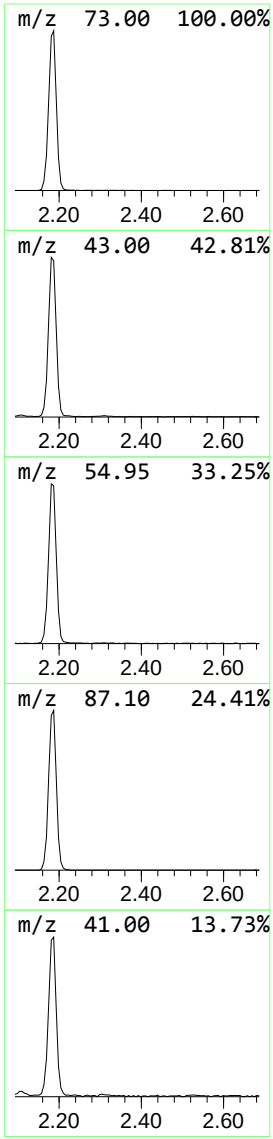
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.187	35.38 ng	1789860	1,4-Dichlorobenzene-d4	6.910

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

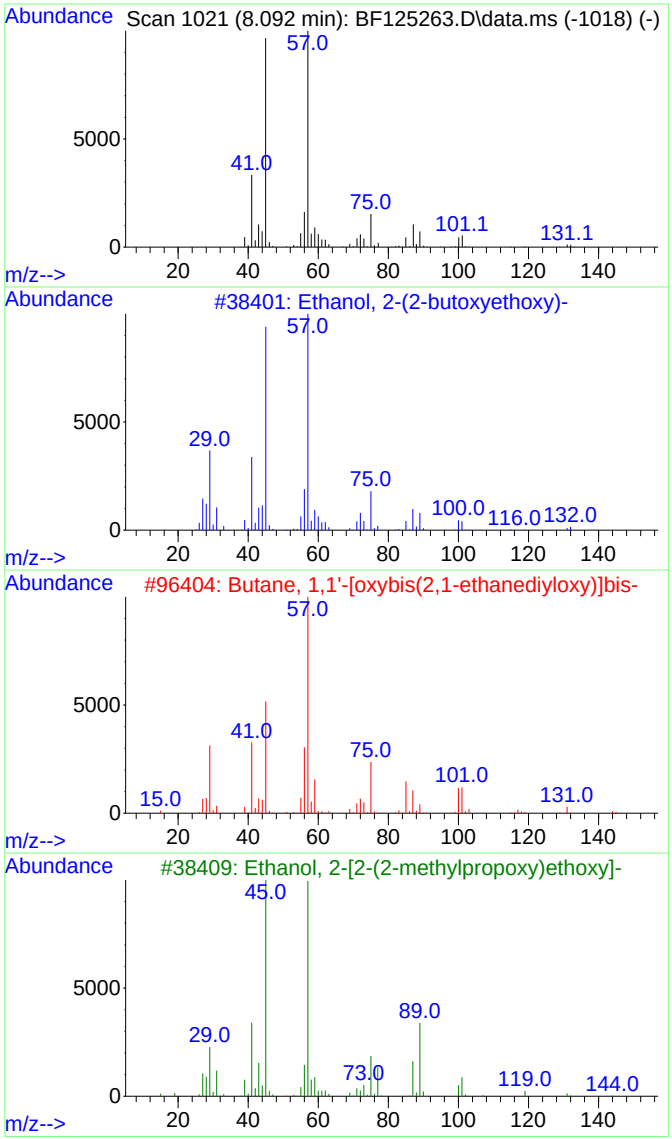
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

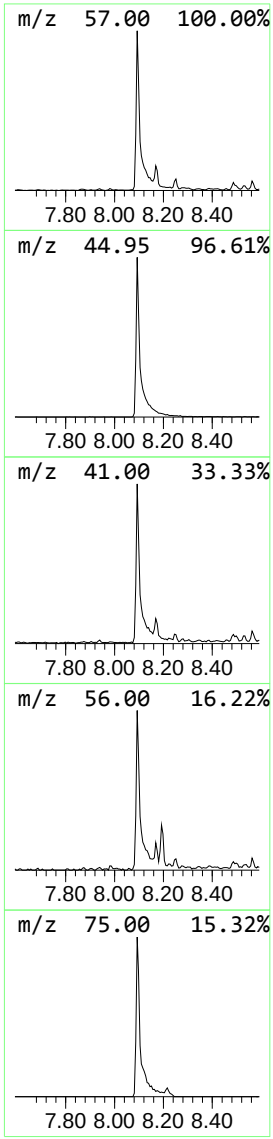
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.092	20.70 ng	1268200	Naphthalene-d8	8.192

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	64
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
4	Tetraethylene glycol	194	C8H18O5	000112-60-7	50
5	Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

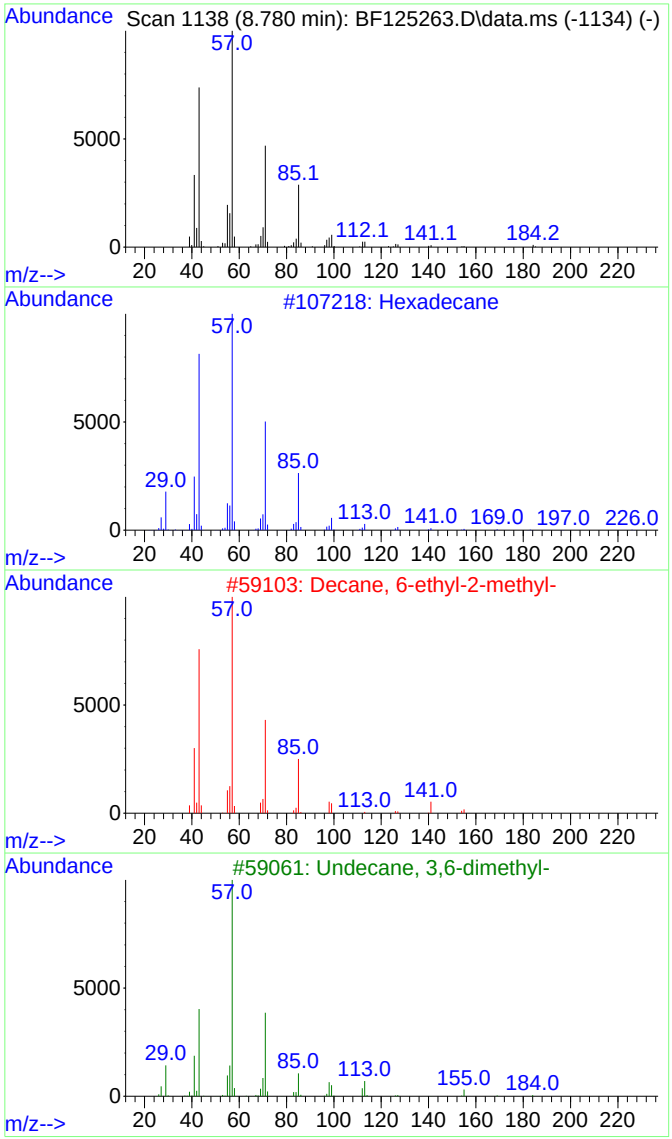
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

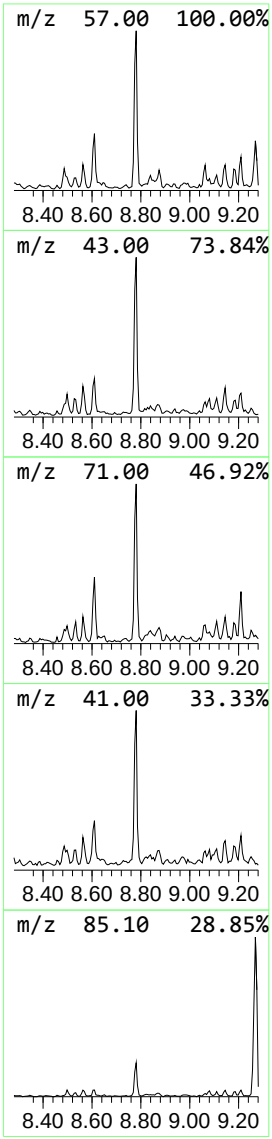
****

Peak Number 3 Hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.780	5.11 ng	313104	Naphthalene-d8	8.192

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	86
2	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	86
3	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	83
4	Tetradecane	198	C14H30	000629-59-4	78
5	Pentadecane	212	C15H32	000629-62-9	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
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Acq On : 30 Aug 2021 15:32
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Sample : M3569-02
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ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

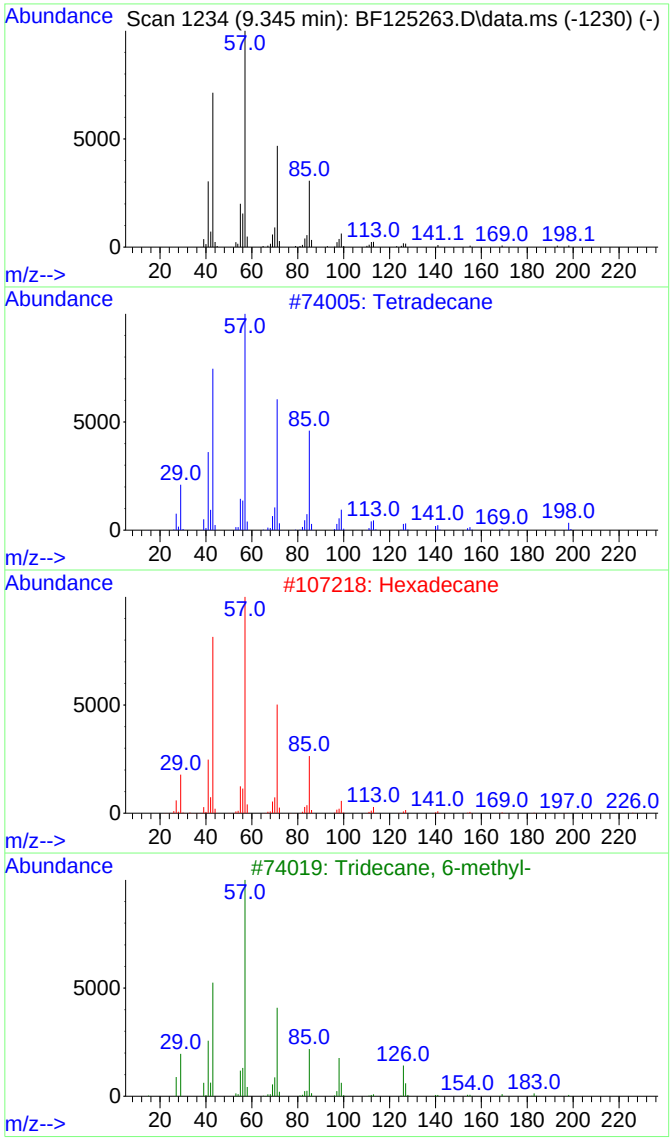
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

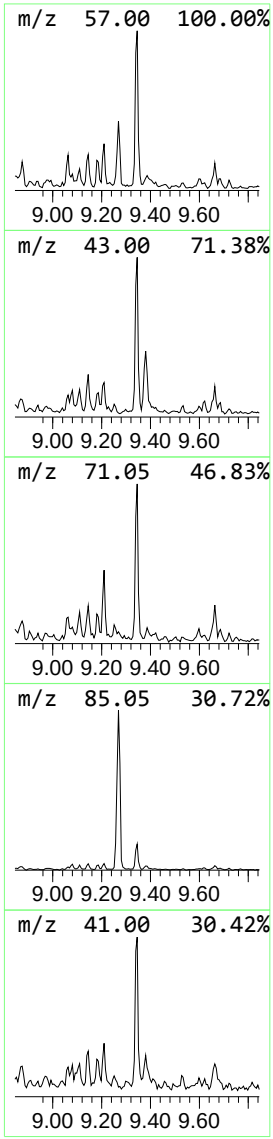
****

Peak Number 4 Tetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.345	2.94 ng	224176	Acenaphthene-d10	9.951

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	94
2	Hexadecane	226	C16H34	000544-76-3	90
3	Tridecane, 6-methyl-	198	C14H30	013287-21-3	87
4	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86
5	Undecane	156	C11H24	001120-21-4	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

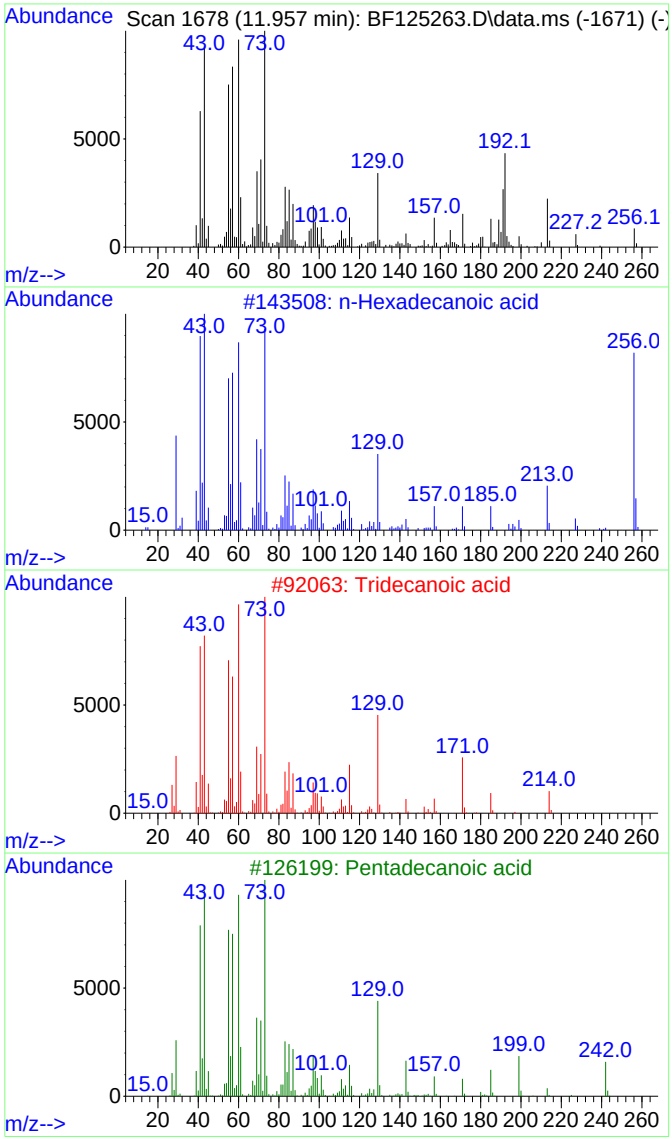
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

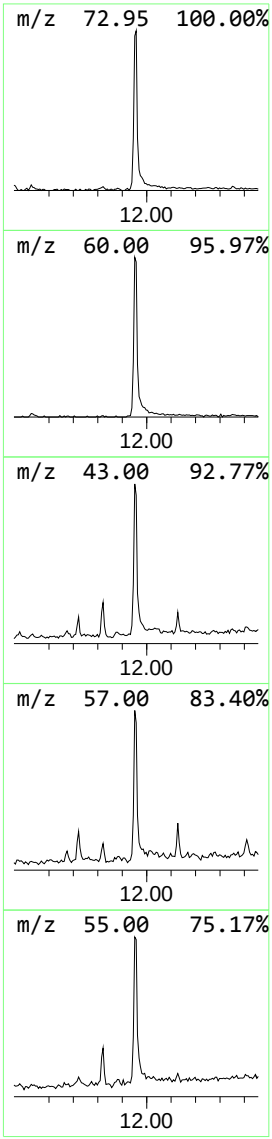
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.957	11.15 ng	979749	Phenanthrene-d10	11.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	78
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	70
4	Undecanoic acid	186	C11H22O2	000112-37-8	62
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S-2

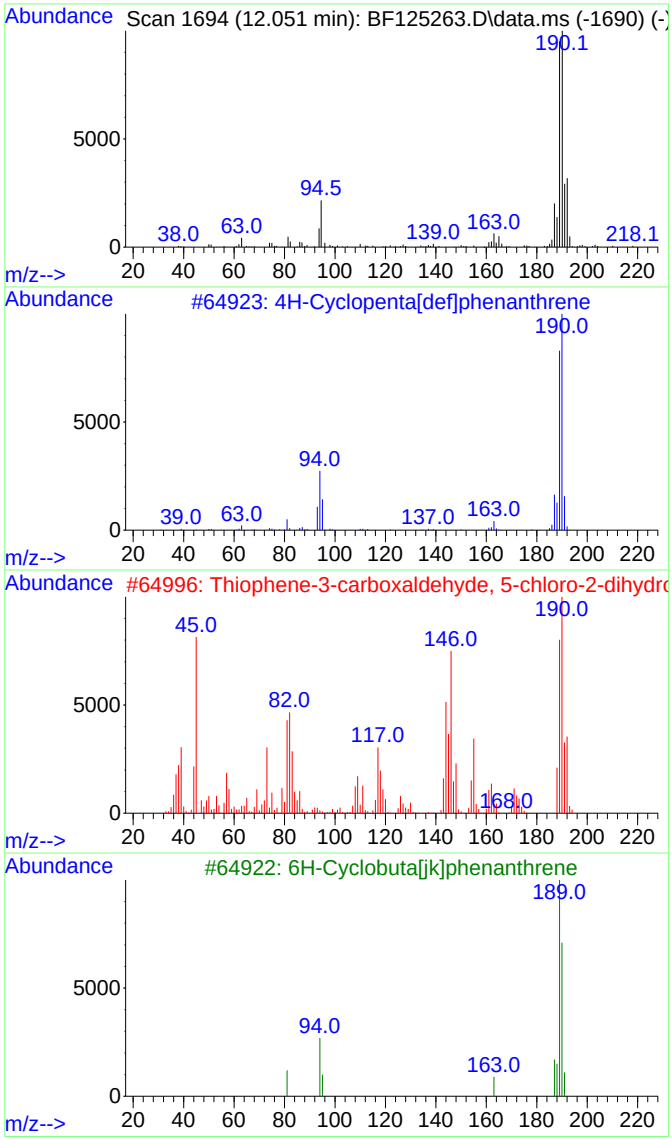
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

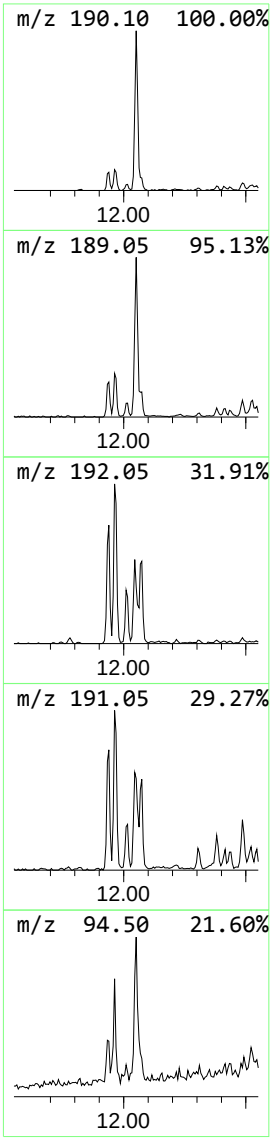
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.051	3.88 ng	341193	Phenanthrene-d10	11.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083021\
Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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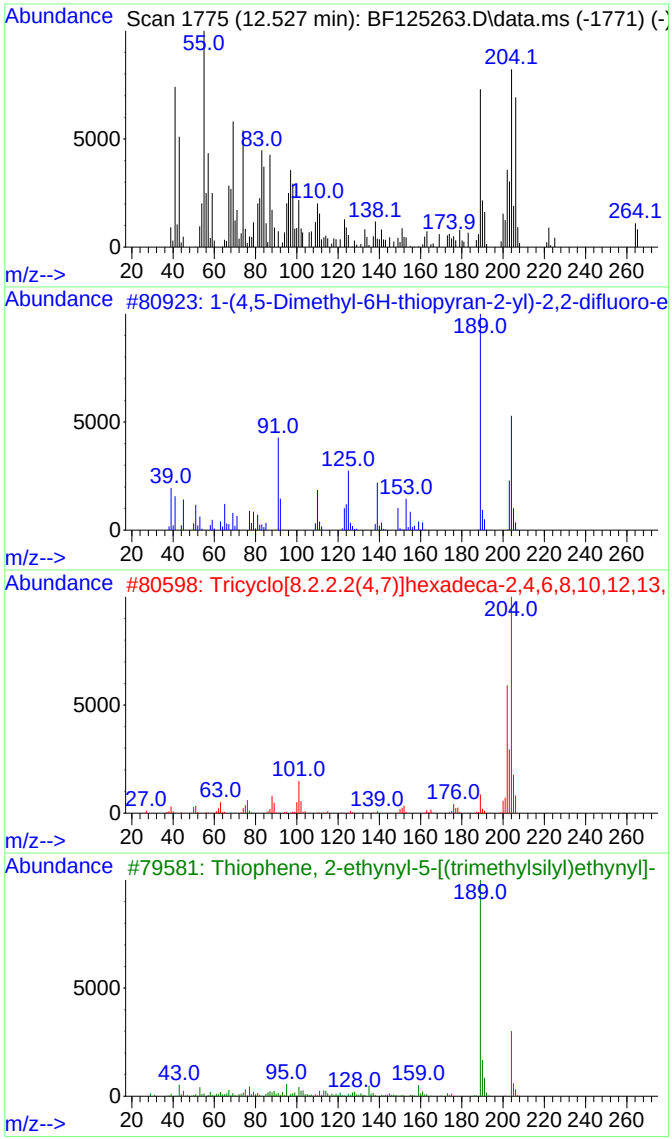
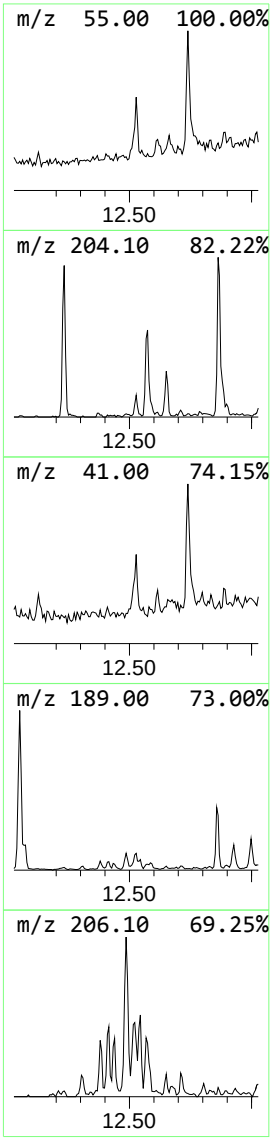
Peak Number 7 unknown12.527 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.527	2.56 ng	224694	Phenanthrene-d10	11.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(4,5-Dimethyl-6H-thiopyran-2-y...	204	C9H10F2OS	1000318-25-0	43
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	38
3	Thiophene, 2-ethynyl-5-[(trimeth...	204	C11H12SSi	182323-29-1	38
4	3,5-Dichloro-2,4-dimethyl-1-meth...	204	C9H10Cl2O	1010250-17-8	38
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	30





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Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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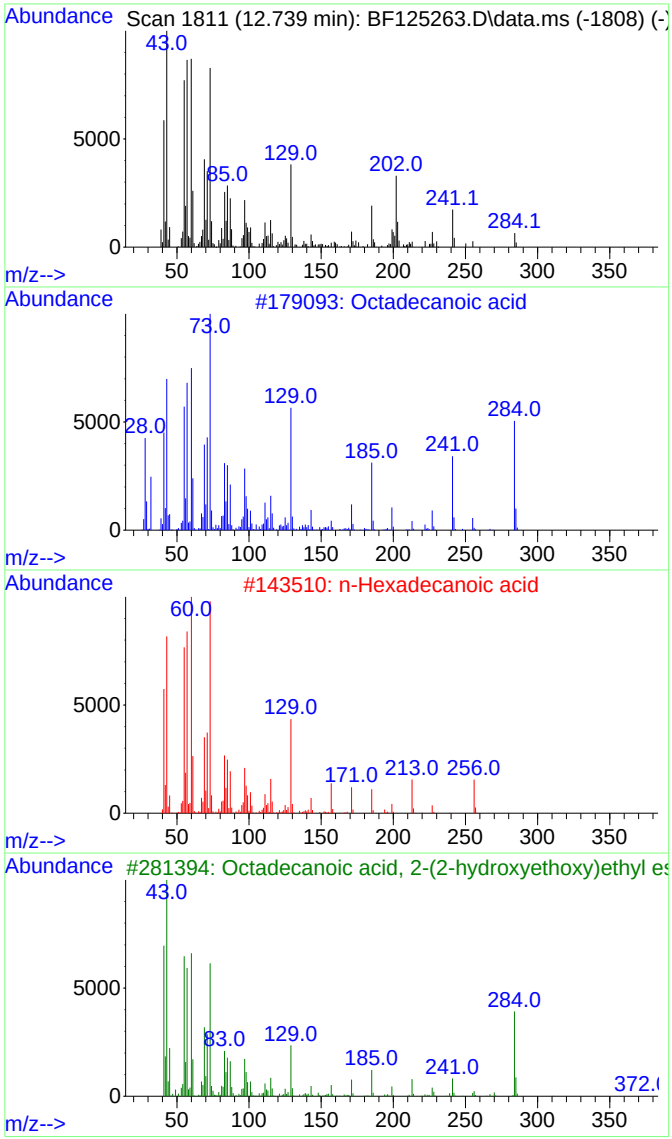
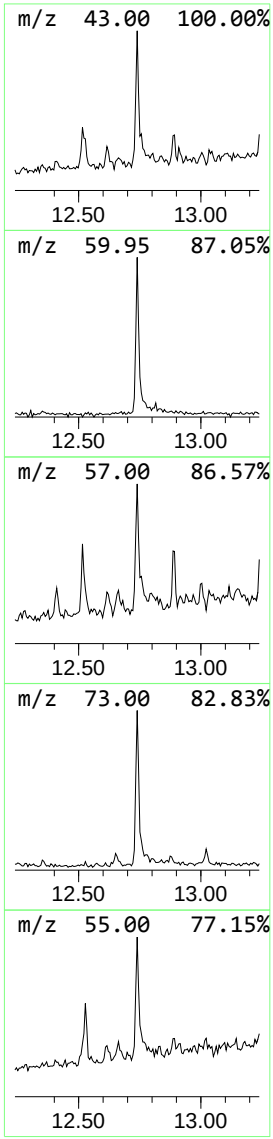
Peak Number 8 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.739	3.17 ng	278701	Phenanthrene-d10	11.439

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
3	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	Tridecanoic acid	214	C13H26O2	000638-53-9	72





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Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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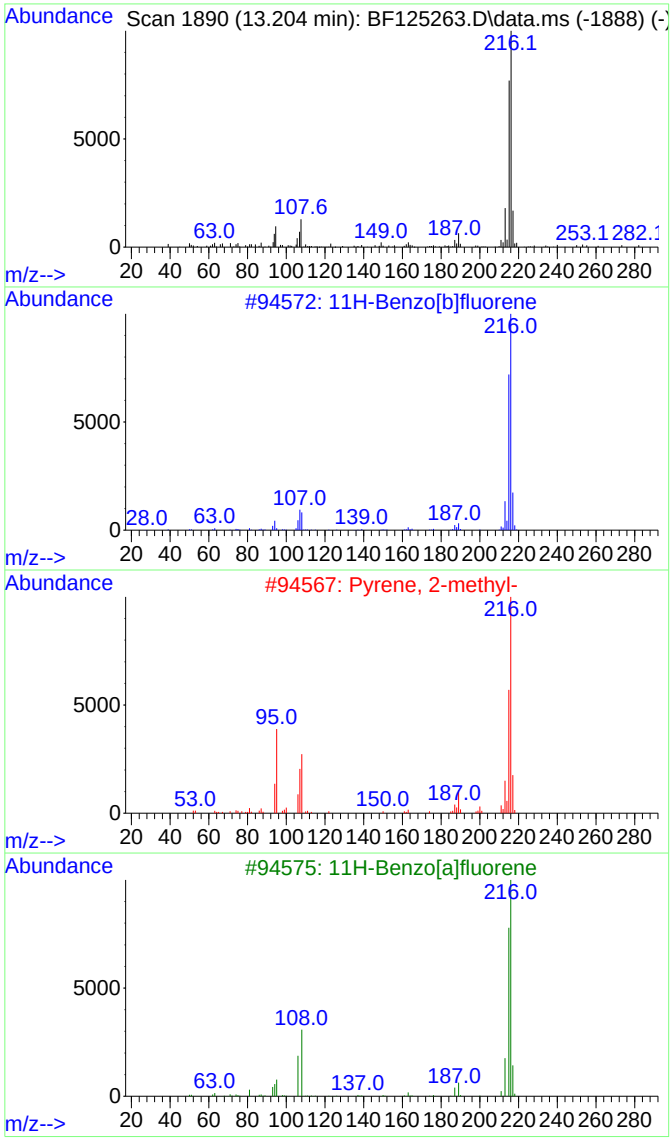
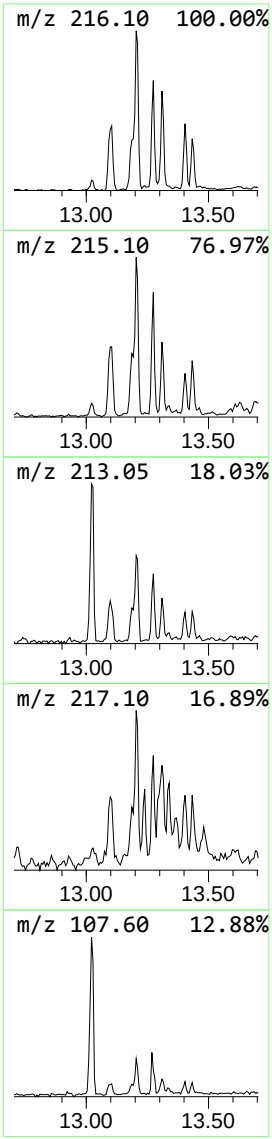
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.204	2.00 ng	213684	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	80





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Data File : BF125263.D
Acq On : 30 Aug 2021 15:32
Operator : JU/CG
Sample : M3569-02
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

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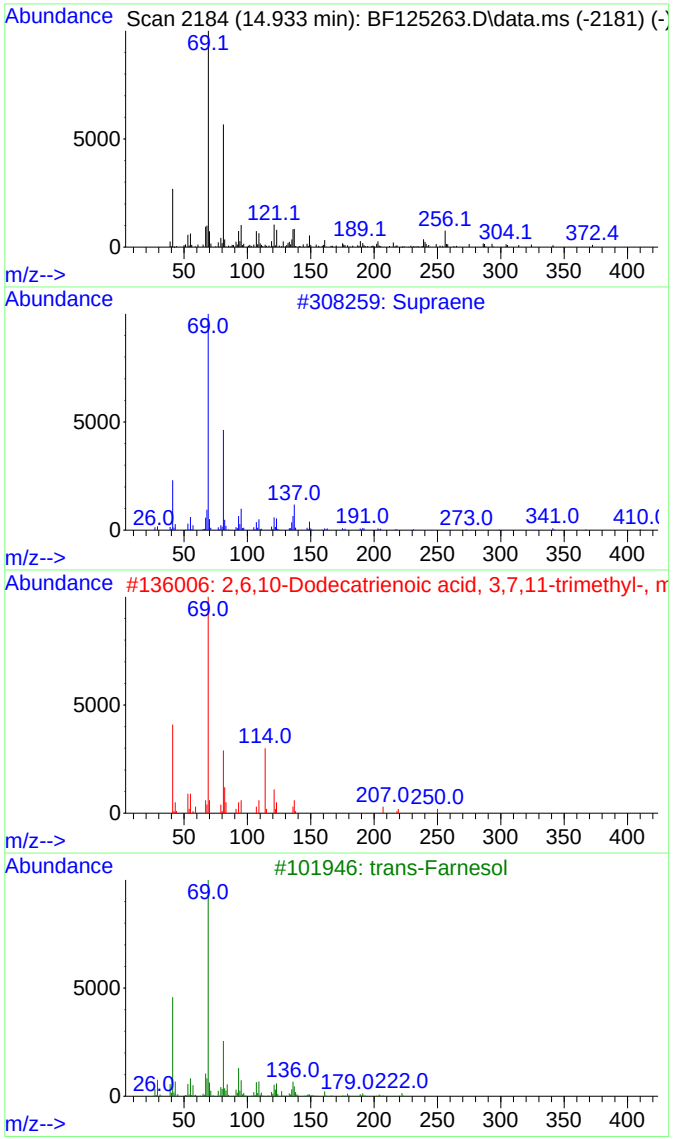
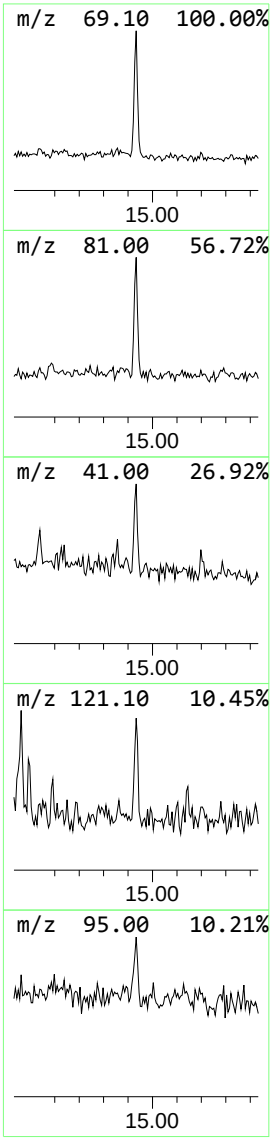
Peak Number 10 Supraene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.933	2.13 ng	131481	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	83
2	2,6,10-Dodecatrienoic acid, 3,7,...	250	C16H26O2	004176-79-8	72
3	trans-Farnesol	222	C15H26O	000106-28-5	72
4	(3E,7E)-4,8,12-Trimethyltrideca-...	218	C16H26	062235-06-7	68
5	1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	64





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Sample : M3569-02
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ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : S-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
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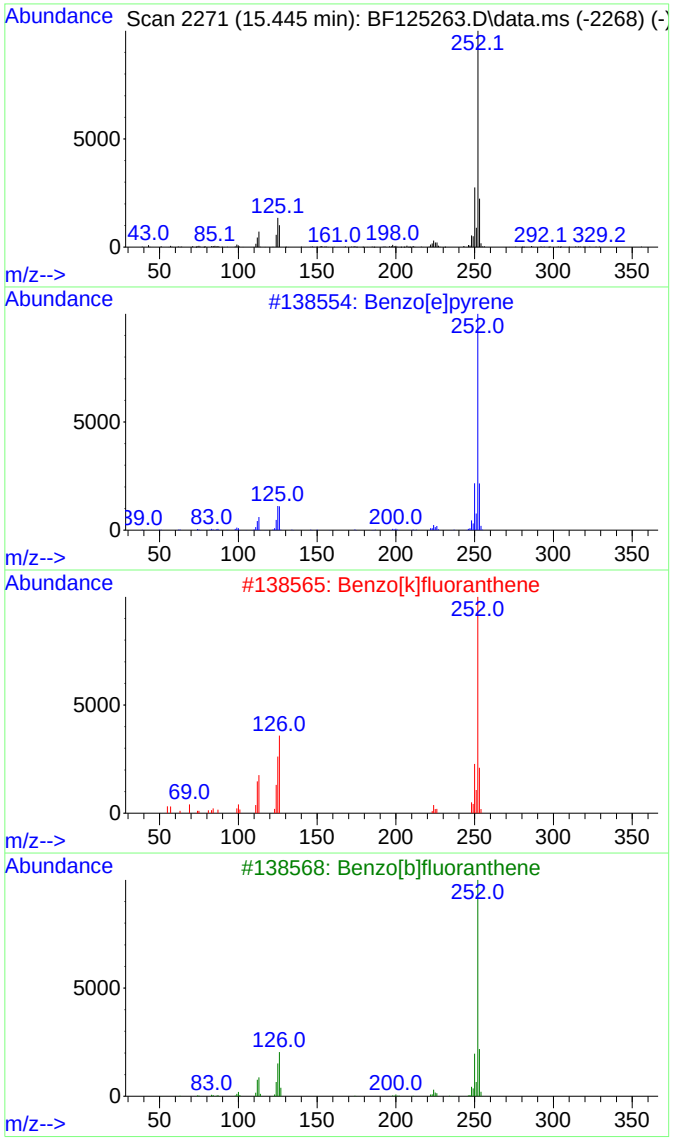
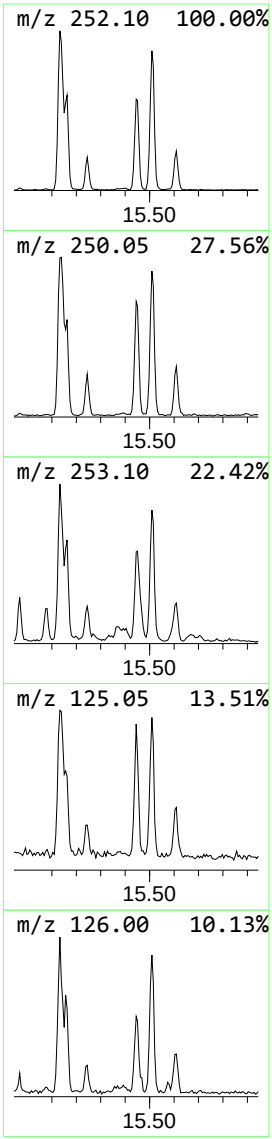
Peak Number 12 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.445	9.22 ng	568017	Perylene-d12	15.580

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4	Benzo[a]pyrene	252	C20H12	000050-32-8	94
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	89





Tentatively Identified Compound (LSC) summary

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ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	2.187	35.4	ng	1789860	1	6.910	1011740	20.0
Ethanol, 2-(2-b...	8.092	20.7	ng	1268200	2	8.192	1225520	20.0
Hexadecane	8.780	5.1	ng	313104	2	8.192	1225520	20.0
Tetradecane	9.345	2.9	ng	224176	3	9.951	1524660	20.0
n-Hexadecanoic ...	11.957	11.2	ng	979749	4	11.439	1757240	20.0
4H-Cyclopenta[d...	12.051	3.9	ng	341193	4	11.439	1757240	20.0
unknown12.527	12.527	2.6	ng	224694	4	11.439	1757240	20.0
Octadecanoic acid	12.739	3.2	ng	278701	4	11.439	1757240	20.0
11H-Benzo[b]flu...	13.204	2.0	ng	213684	5	14.080	2131900	20.0
Supraene	14.933	2.1	ng	131481	6	15.580	1232200	20.0
Benzo[e]pyrene	15.445	9.2	ng	568017	6	15.580	1232200	20.0