

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122121\
 Data File : BF126318.D
 Acq On : 22 Dec 2021 2:29
 Operator : JU/CG
 Sample : M5184-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-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-BF121521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF126318.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	7	18	rVB	123232	159089	3.66%	0.453%
2	5.040	493	502	513	rBV	63536	87623	2.02%	0.250%
3	5.434	563	569	572	rBV	3759592	3843488	88.51%	10.949%
4	6.445	735	741	744	rBV	3463071	4012476	92.41%	11.430%
5	6.816	800	804	807	rBV	1395798	1170428	26.95%	3.334%
6	7.375	893	899	902	rBV	2542073	2740528	63.11%	7.807%
7	8.098	1017	1022	1027	rBV	1768863	1634425	37.64%	4.656%
8	9.175	1200	1205	1217	rBV	4030867	4342212	100.00%	12.369%
9	9.257	1217	1219	1229	rVB3	57721	82858	1.91%	0.236%
10	9.851	1315	1320	1337	rBV	1741658	1800064	41.46%	5.128%
11	10.639	1449	1454	1465	rBV	2414946	2618740	60.31%	7.460%
12	11.339	1569	1573	1575	rBV	1722693	1704143	39.25%	4.854%
13	11.357	1575	1576	1579	rVB	149874	100185	2.31%	0.285%
14	11.875	1660	1664	1667	rBV2	399545	432041	9.95%	1.231%
15	11.904	1667	1669	1673	rVB	61042	63368	1.46%	0.181%
16	12.545	1775	1778	1781	rBV	437626	373145	8.59%	1.063%
17	12.657	1793	1797	1802	rBV2	88599	113920	2.62%	0.325%
18	12.774	1810	1817	1821	rVV	374982	395169	9.10%	1.126%
19	12.827	1823	1826	1830	rVB3	53214	49025	1.13%	0.140%
20	12.927	1838	1843	1846	rBV	3841620	3676752	84.67%	10.474%
21	13.104	1871	1873	1877	rVB	58747	48246	1.11%	0.137%
22	13.204	1887	1890	1893	rBV	45617	49098	1.13%	0.140%
23	13.610	1956	1959	1964	rVB2	43357	58887	1.36%	0.168%
24	13.816	1991	1994	1999	rVB3	38548	51683	1.19%	0.147%
25	13.898	2002	2008	2016	rBV3	323921	1096192	25.25%	3.123%
26	13.974	2016	2021	2023	rVV2	1367373	1610841	37.10%	4.589%
27	13.998	2023	2025	2032	rVV	262230	291536	6.71%	0.830%
28	14.068	2032	2037	2045	rVB	391724	468236	10.78%	1.334%
29	14.757	2148	2154	2160	rVB4	32202	58235	1.34%	0.166%
30	14.833	2163	2167	2176	rVB2	40172	55576	1.28%	0.158%
31	15.004	2190	2196	2199	rBV	164083	271856	6.26%	0.774%
32	15.092	2207	2211	2219	rVB3	40188	70404	1.62%	0.201%
33	15.298	2241	2246	2251	rVB	96518	120650	2.78%	0.344%
34	15.357	2252	2256	2262	rVV	107779	136627	3.15%	0.389%
35	15.421	2262	2267	2279	rVV	712890	1004382	23.13%	2.861%
36	15.898	2344	2348	2352	rVB2	29925	43631	1.00%	0.124%

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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-BF121521.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.139	2377	2389	2390	rBV9	28768	87315	2.01%	0.249%
38	16.851	2504	2510	2520	rVB2	44060	102790	2.37%	0.293%
39	17.280	2577	2583	2589	rBV2	34399	79127	1.82%	0.225%

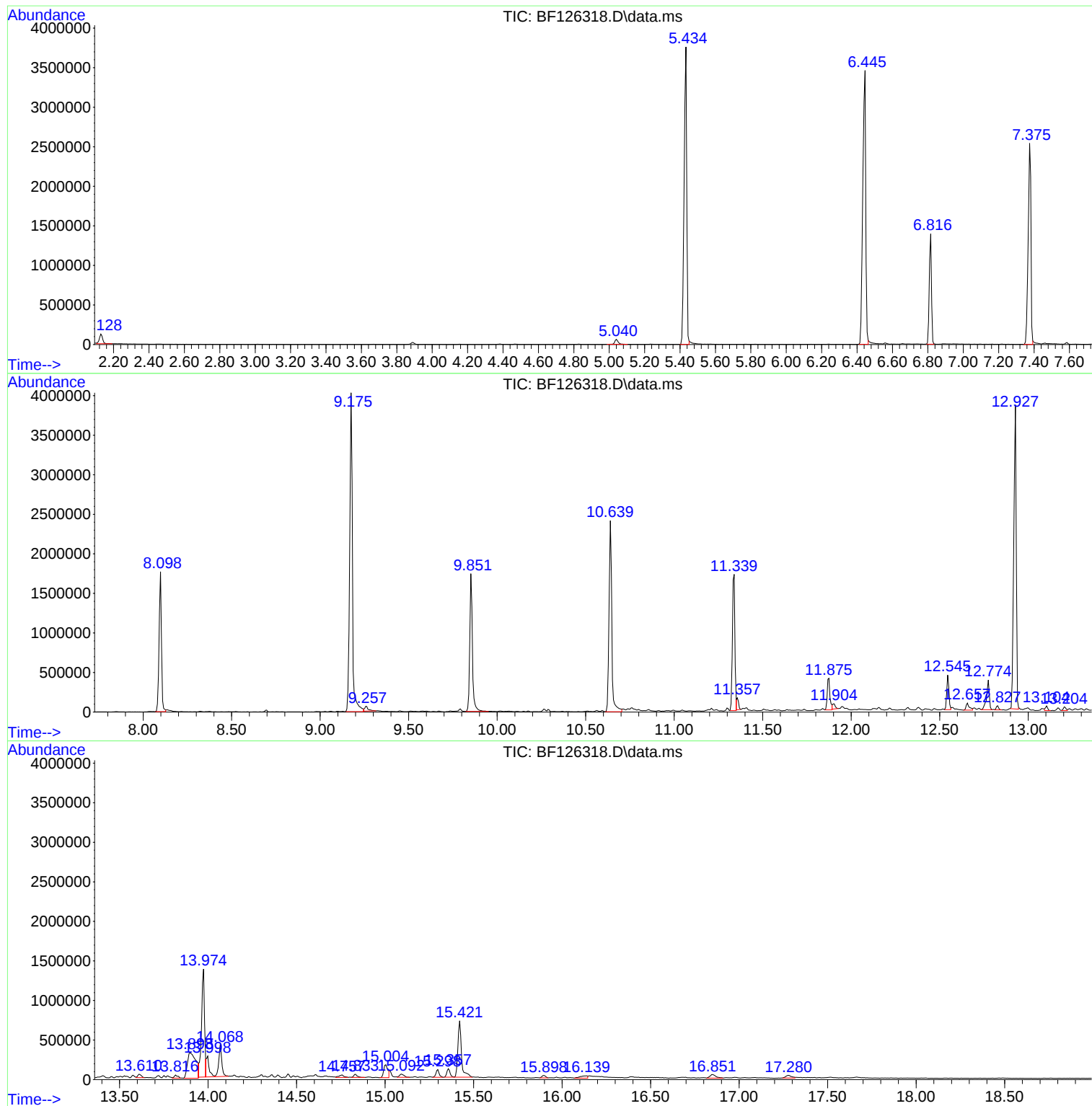
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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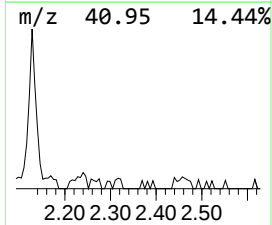
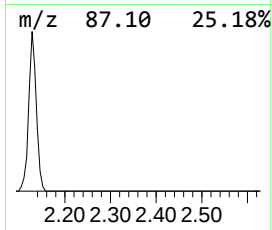
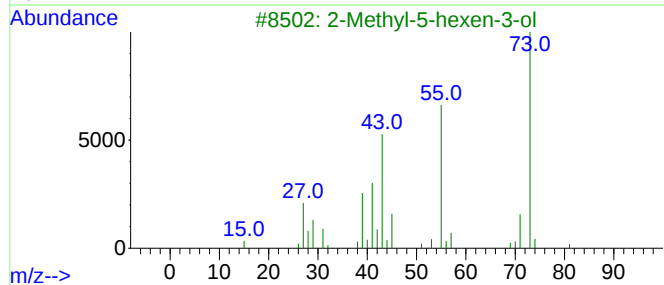
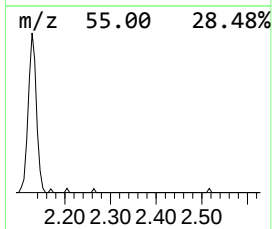
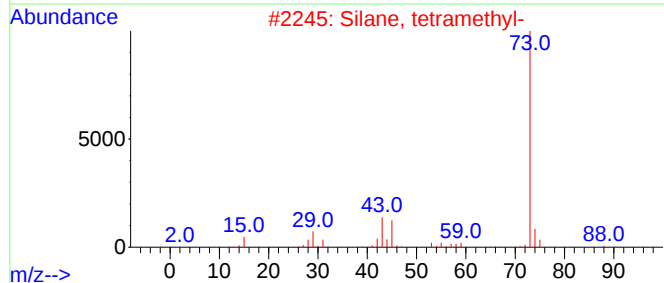
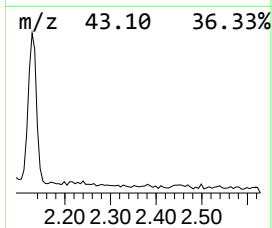
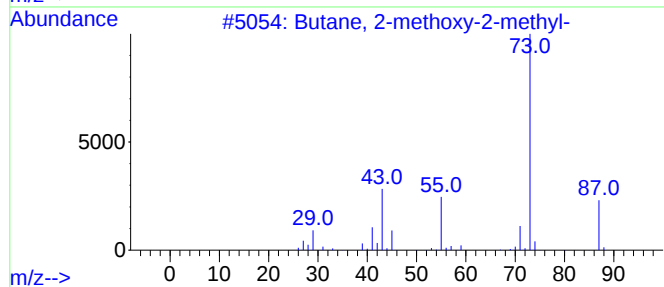
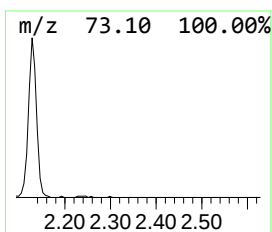
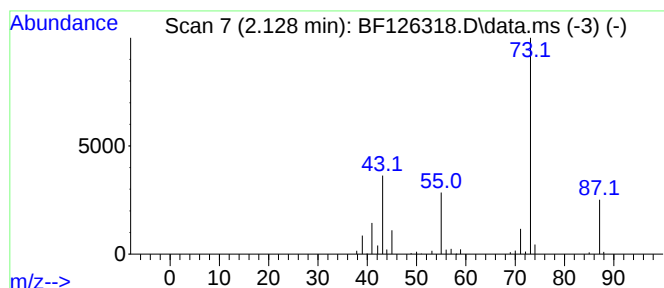
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.128	2.72 ng	159089	1,4-Dichlorobenzene-d4	6.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	28
4			3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	16
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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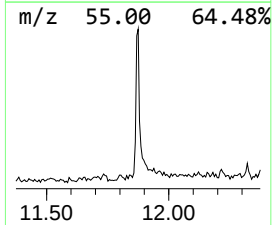
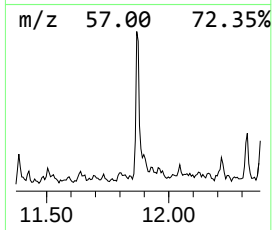
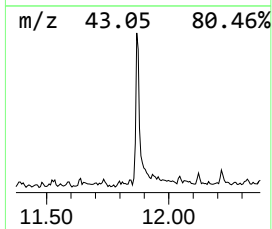
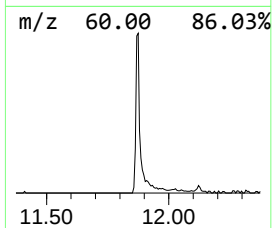
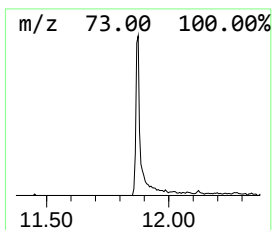
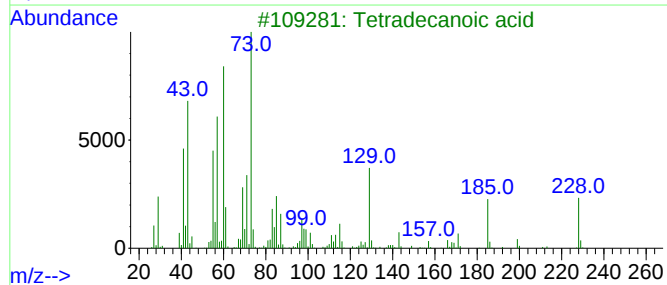
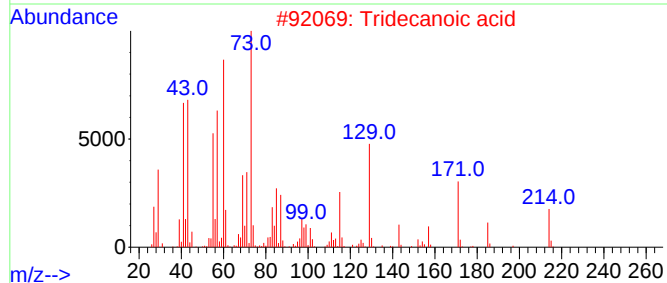
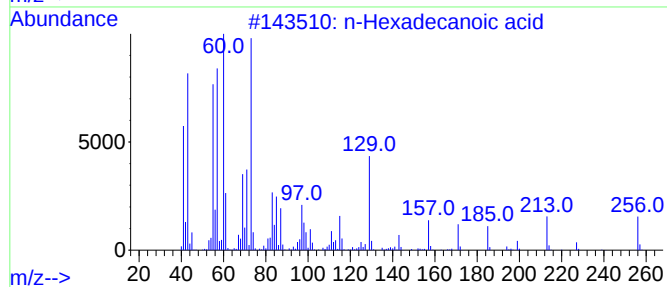
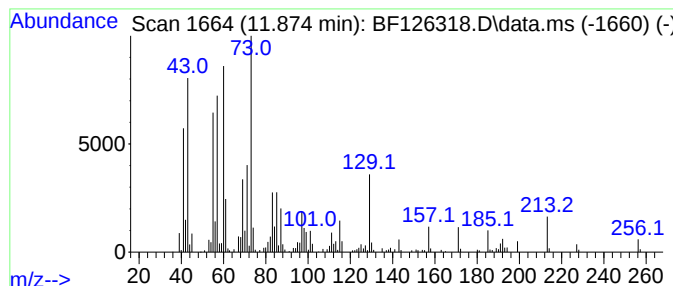
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.874	5.07 ng	432041	Phenanthrene-d10	11.339

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	94
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	64



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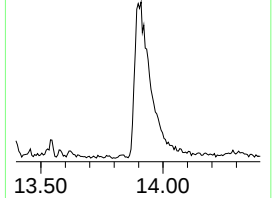
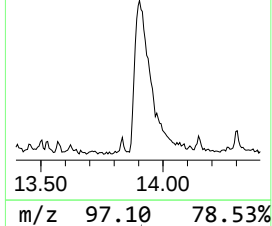
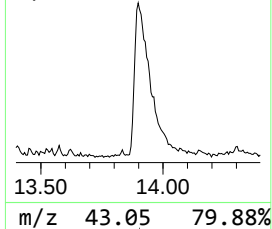
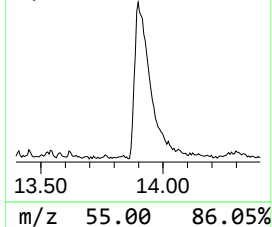
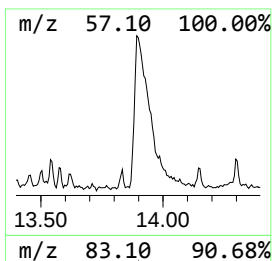
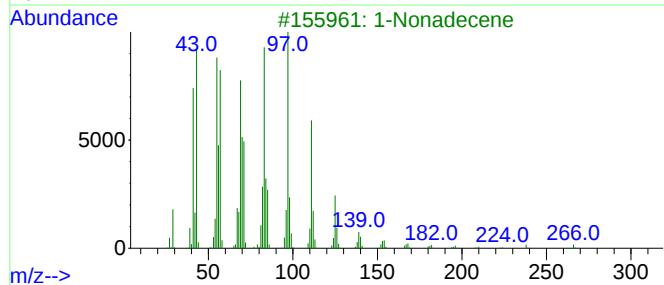
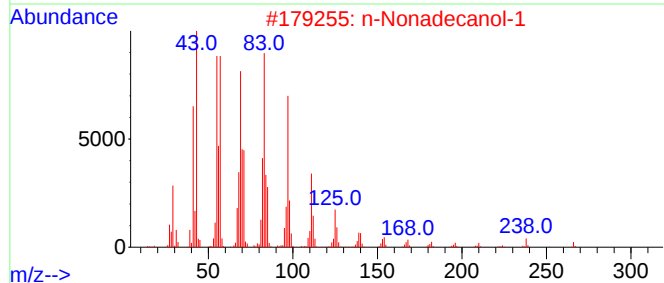
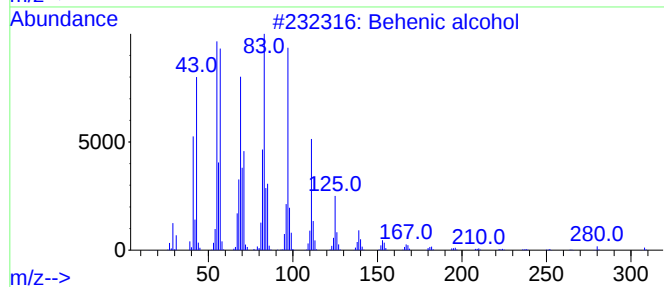
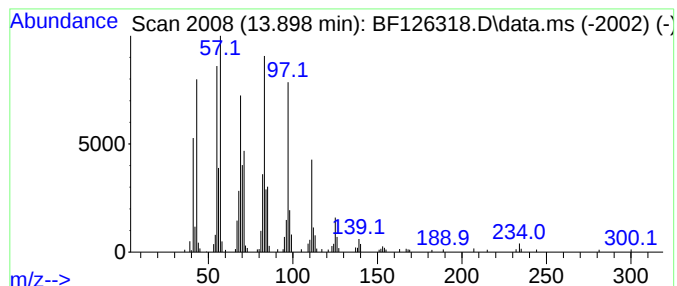
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Behenic alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.898	13.61 ng	1096190	Chrysene-d12	13.974

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	94
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	93
3			1-Nonadecene	266	C19H38	018435-45-5	91
4			5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



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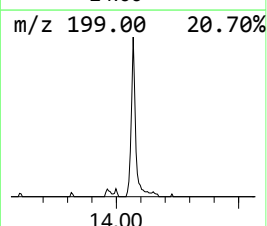
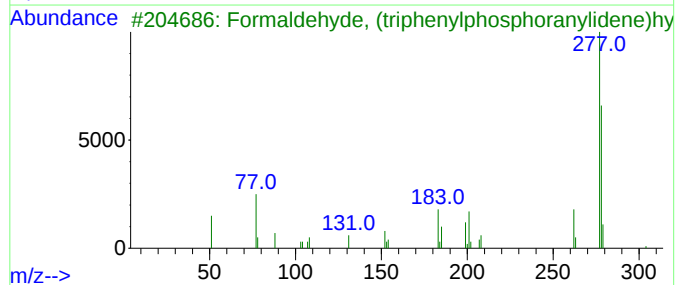
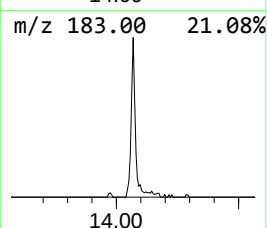
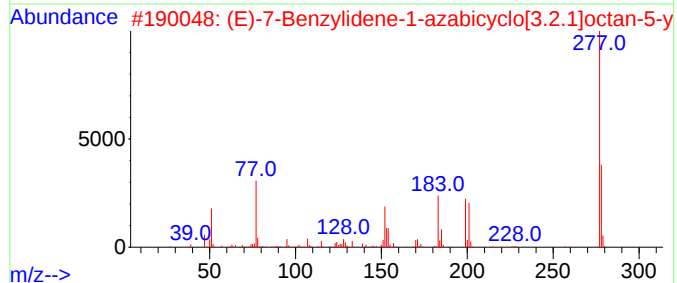
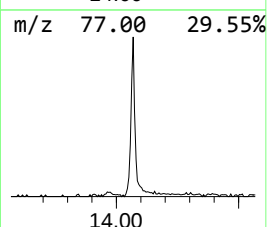
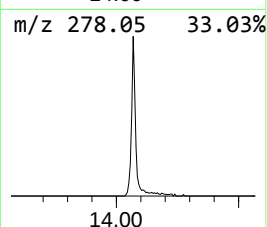
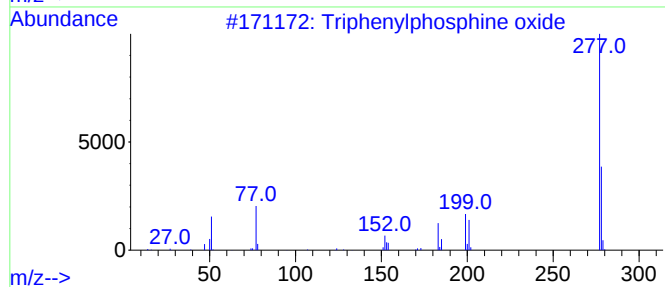
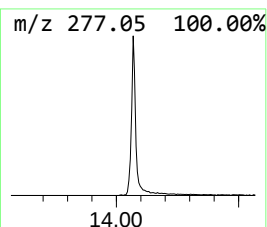
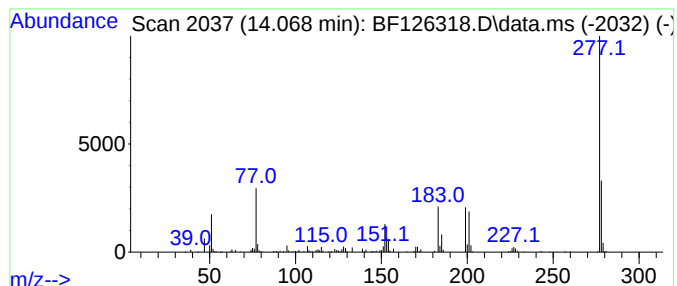
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Triphenylphosphine oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.069	5.81 ng	468236	Chrysene-d12	13.974

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19NO3S	1000483-83-4	94
3			Formaldehyde, (triphenylphosphor...]	304	C19H17N2P	015990-54-2	64
4			4-Nitro-4'-methyldiphenylsulfone	277	C13H11NO4S	004094-37-5	46
5			2-Phenyl-6-(trifluoromethyl)-1H-...	277	C15H10F3NO	1000387-59-3	37



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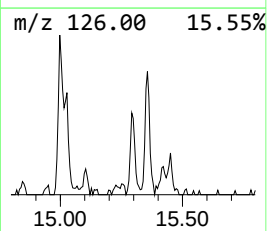
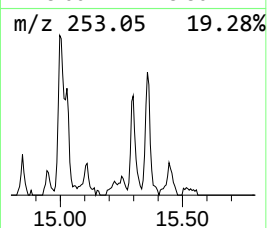
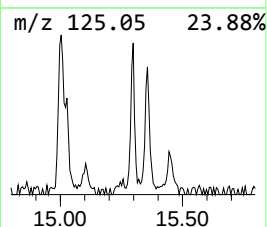
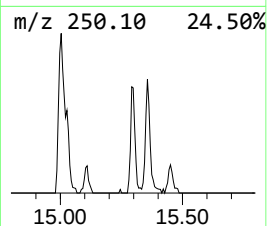
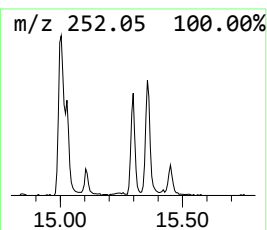
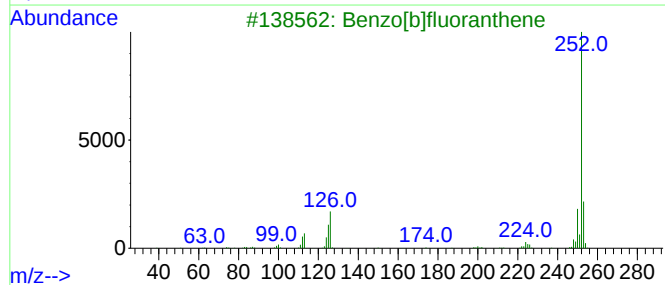
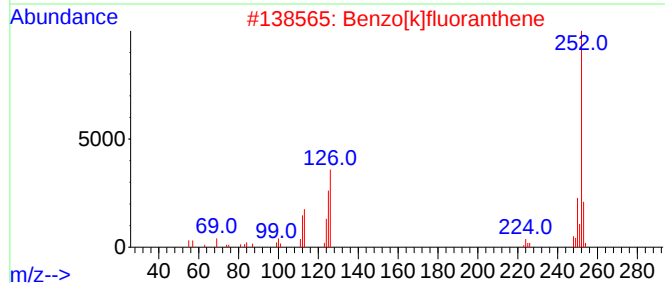
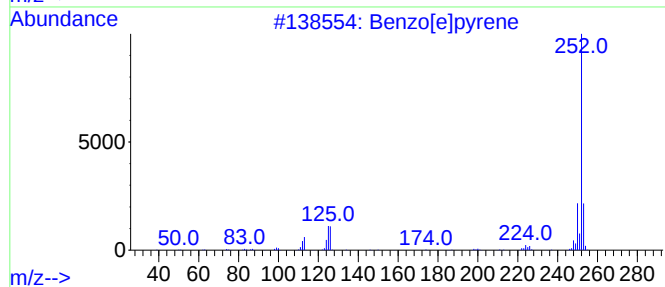
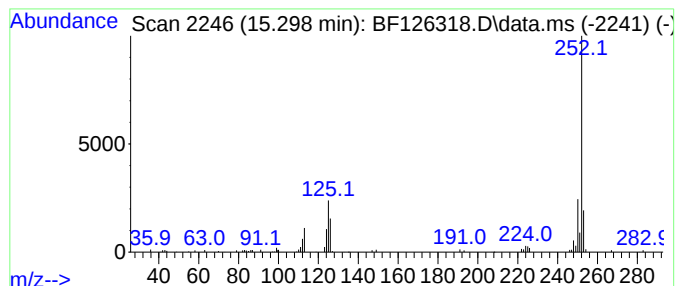
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.298	2.40 ng	120650	Perylene-d12	15.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	95
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Perylene	252	C20H12	000198-55-0	90
5			Benzo[a]pyrene	252	C20H12	000050-32-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122121\
Data File : BF126318.D
Acq On : 22 Dec 2021 2:29
Operator : JU/CG
Sample : M5184-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.128	2.7	ng	159089	1	6.816	1170430	20.0
n-Hexadecanoic ...	11.874	5.1	ng	432041	4	11.339	1704140	20.0
Behenic alcohol	13.898	13.6	ng	1096190	5	13.974	1610840	20.0
Triphenylphosph...	14.069	5.8	ng	468236	5	13.974	1610840	20.0
Benzo[e]pyrene	15.298	2.4	ng	120650	6	15.421	1004380	20.0