

Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
 Data File : BF097900.D
 Acq On : 21 Aug 2017 12:18
 Operator : SJ/JU
 Sample : I4869-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-TP-106-A(0-5)COMP

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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	3.134	175	178	188	rBV	50890	101539	2.47%	0.267%
2	4.951	483	487	496	rVB	294665	457701	11.14%	1.202%
3	5.363	551	557	560	rBV	3388900	3804737	92.63%	9.994%
4	6.387	725	731	734	rBV	3304323	4107665	100.00%	10.789%
5	6.522	748	754	757	rBV	3408399	3749597	91.28%	9.849%
6	6.745	789	792	796	rBV	985123	803573	19.56%	2.111%
7	6.904	815	819	822	rBV	3053841	2770553	67.45%	7.277%
8	7.316	883	889	891	rBV	2239967	2543312	61.92%	6.680%
9	8.028	1006	1010	1014	rBV	1248915	1048172	25.52%	2.753%
10	9.110	1188	1194	1197	rBV	3614445	3553017	86.50%	9.333%
11	9.498	1256	1260	1263	rBV	520517	404216	9.84%	1.062%
12	9.781	1301	1308	1311	rBV2	1228412	1064452	25.91%	2.796%
13	10.222	1379	1383	1386	rVB	50818	47782	1.16%	0.126%
14	10.322	1398	1400	1407	rVB2	44585	45290	1.10%	0.119%
15	10.569	1436	1442	1445	rBV	2198239	2146458	52.25%	5.638%
16	10.692	1456	1463	1470	rBV3	87046	124650	3.03%	0.327%
17	11.263	1556	1560	1562	rBV	1010356	956426	23.28%	2.512%
18	11.286	1562	1564	1567	rVV	468354	371304	9.04%	0.975%
19	11.333	1567	1572	1575	rVB	182593	160057	3.90%	0.420%
20	11.763	1641	1645	1647	rBV	89967	84191	2.05%	0.221%
21	11.786	1647	1649	1653	rVB	140953	122357	2.98%	0.321%
22	11.833	1653	1657	1660	rBV	75092	64983	1.58%	0.171%
23	11.869	1660	1663	1666	rVV	148564	164414	4.00%	0.432%
24	11.892	1666	1667	1673	rVB	67833	45384	1.10%	0.119%
25	12.057	1692	1695	1697	rBV	65621	56120	1.37%	0.147%
26	12.310	1734	1738	1741	rBV	72052	79566	1.94%	0.209%
27	12.345	1741	1744	1749	rVV4	43470	74751	1.82%	0.196%
28	12.392	1749	1752	1756	rVB	56339	64453	1.57%	0.169%
29	12.469	1761	1765	1769	rVV	893813	730605	17.79%	1.919%
30	12.698	1798	1804	1807	rBV2	666674	749115	18.24%	1.968%
31	12.745	1809	1812	1817	rVB2	44933	63655	1.55%	0.167%
32	12.816	1820	1824	1825	rBV	59039	47099	1.15%	0.124%
33	12.845	1825	1829	1833	rVV	2415273	2314147	56.34%	6.078%
34	12.916	1836	1841	1844	rBV3	64874	102439	2.49%	0.269%

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35	12.998	1852	1855	1857	rBV3	49700	56234	1.37%	0.148%
36	13.021	1857	1859	1863	rVB	117607	106329	2.59%	0.279%
37	13.092	1868	1871	1873	rBV	78352	72761	1.77%	0.191%
38	13.127	1873	1877	1880	rBV2	100092	96837	2.36%	0.254%
39	13.251	1895	1898	1901	rBV2	45134	49338	1.20%	0.130%
40	13.421	1921	1927	1930	rBV7	29673	56809	1.38%	0.149%
41	13.527	1938	1945	1950	rBV	73179	117440	2.86%	0.308%
42	13.639	1960	1964	1967	rBV2	66742	90472	2.20%	0.238%
43	13.669	1967	1969	1971	rVV	74454	65377	1.59%	0.172%
44	13.692	1971	1973	1977	rVV2	49811	76478	1.86%	0.201%
45	13.745	1977	1982	1987	rVB2	136503	208945	5.09%	0.549%
46	13.892	1999	2007	2009	rBV2	911644	1248935	30.40%	3.281%
47	13.916	2009	2011	2018	rVB	350428	324326	7.90%	0.852%
48	14.068	2035	2037	2041	rVB2	37463	43860	1.07%	0.115%
49	14.274	2068	2072	2075	rBV	94245	84754	2.06%	0.223%
50	14.310	2075	2078	2082	rVB	55471	50089	1.22%	0.132%
51	14.368	2085	2088	2091	rBV3	42689	41199	1.00%	0.108%
52	14.739	2148	2151	2157	rBV3	66572	88661	2.16%	0.233%
53	14.904	2174	2179	2182	rBV	291665	441730	10.75%	1.160%
54	15.004	2193	2196	2202	rVB	84120	107562	2.62%	0.283%
55	15.186	2224	2227	2233	rVB	169020	210989	5.14%	0.554%
56	15.245	2233	2237	2243	rVB	272360	318575	7.76%	0.837%
57	15.310	2243	2248	2251	rBV	570553	705390	17.17%	1.853%
58	15.762	2322	2325	2329	rBV4	31372	45183	1.10%	0.119%
59	16.674	2475	2480	2488	rVB2	85961	167004	4.07%	0.439%
60	17.086	2544	2550	2561	rBV	71636	172303	4.19%	0.453%

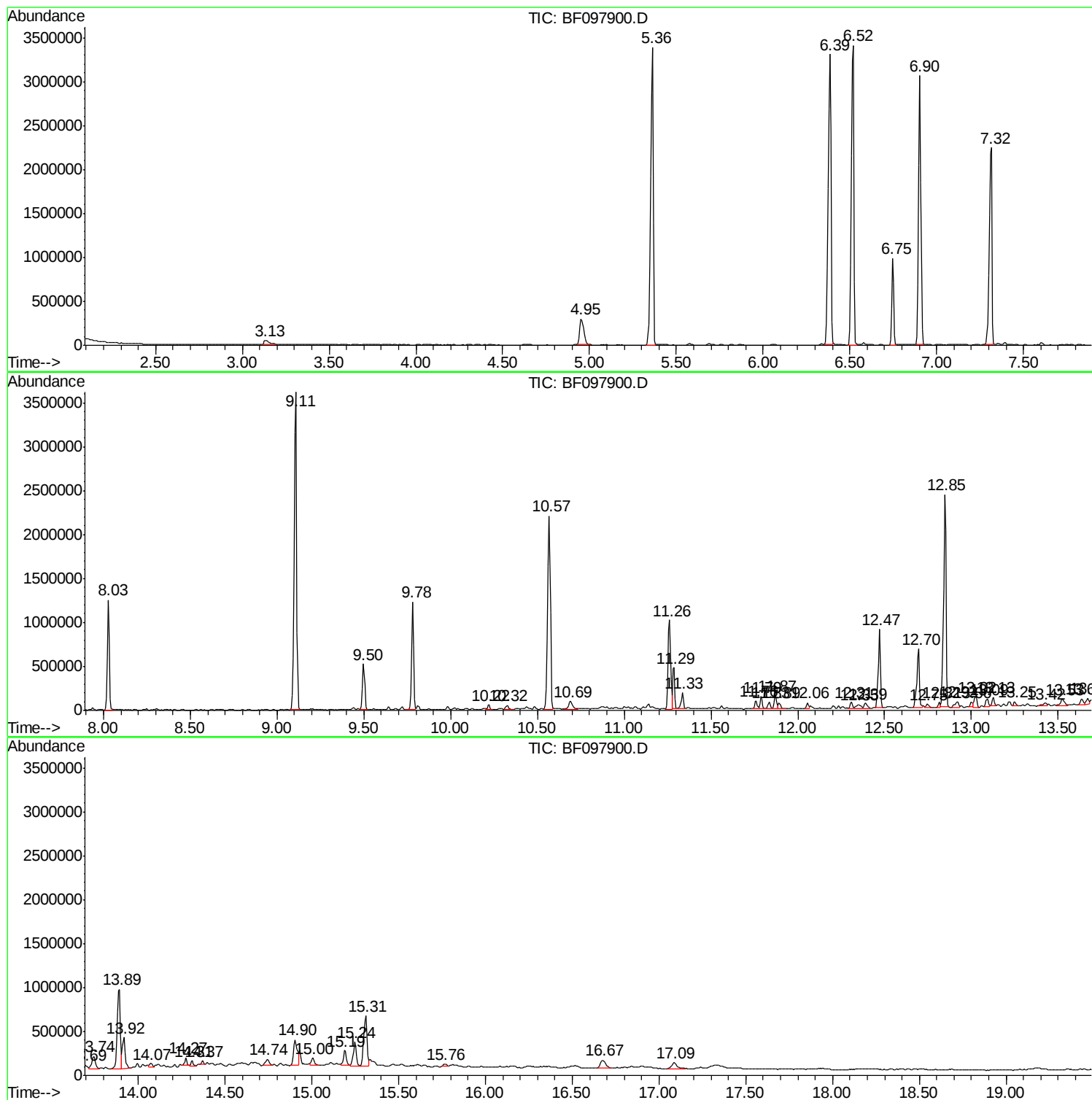
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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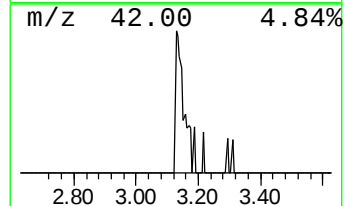
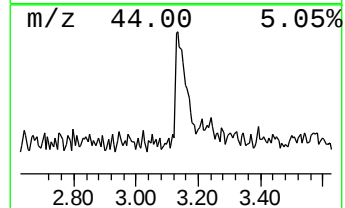
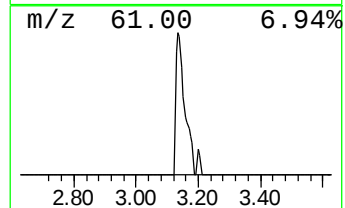
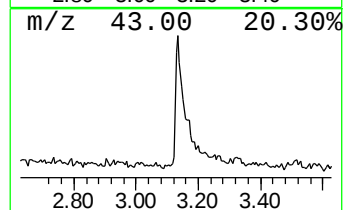
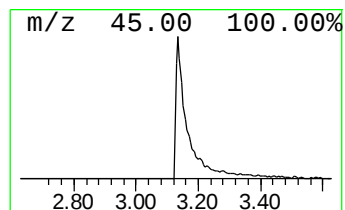
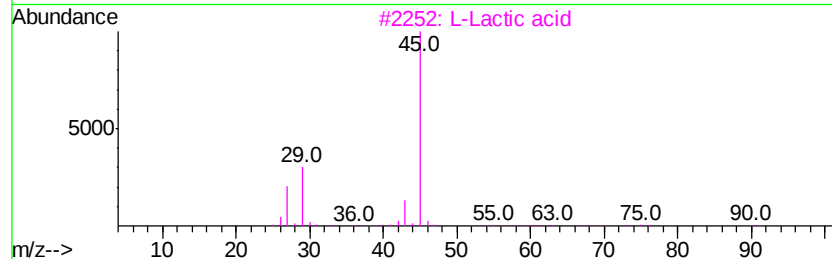
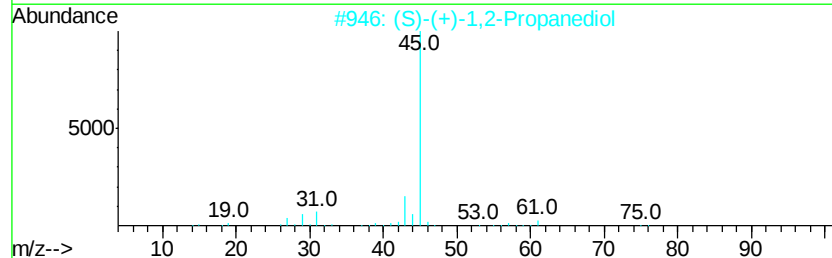
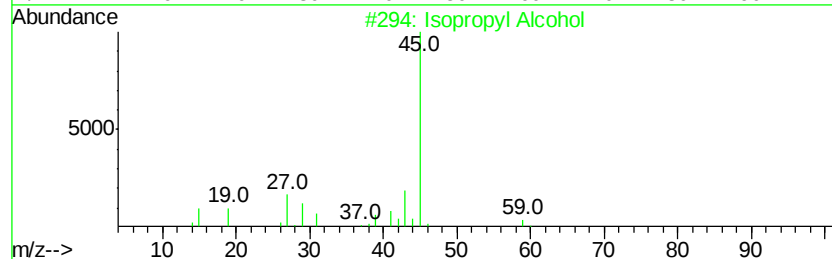
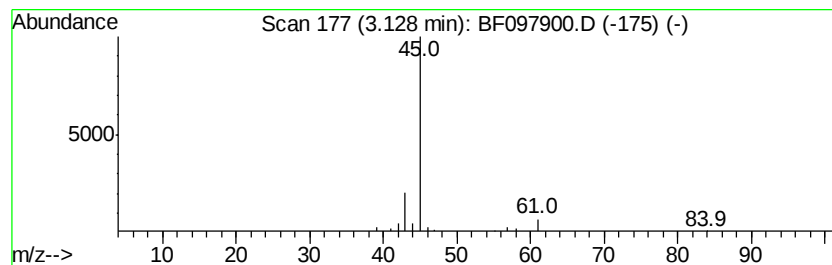
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.13 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	2.53 ng	101539	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isopropyl Alcohol	60	C3H8O	000067-63-0	9
2		(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
3		L-Lactic acid	90	C3H6O3	000079-33-4	9
4		R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9
5		Propylene Glycol	76	C3H8O2	000057-55-6	9



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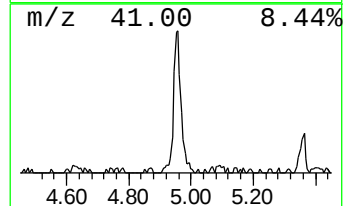
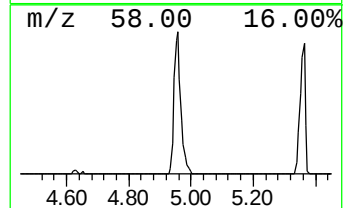
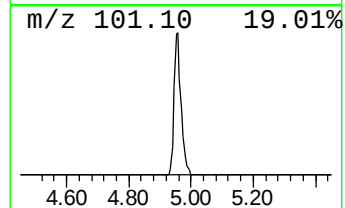
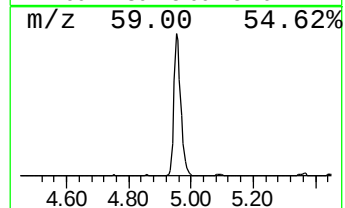
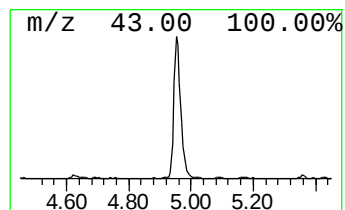
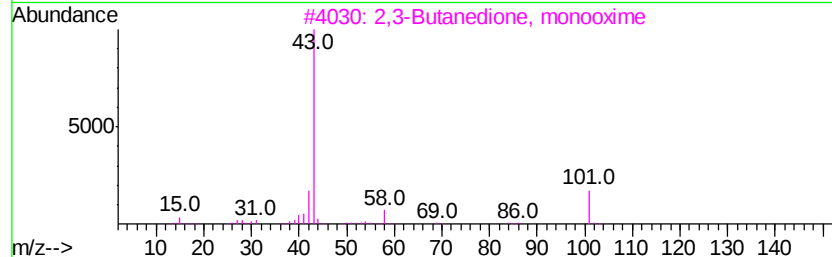
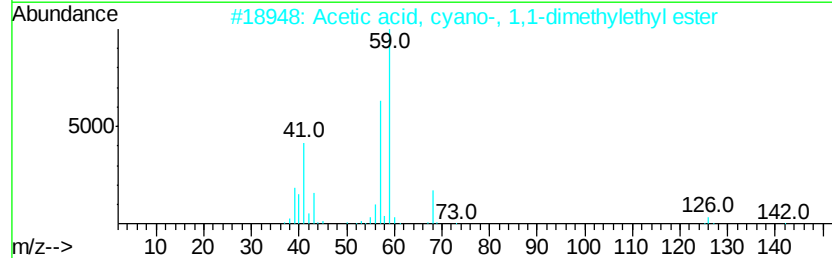
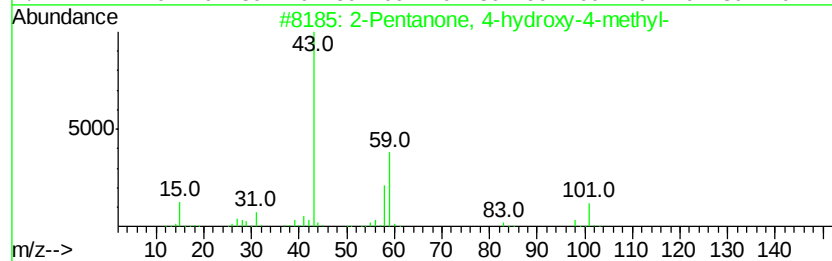
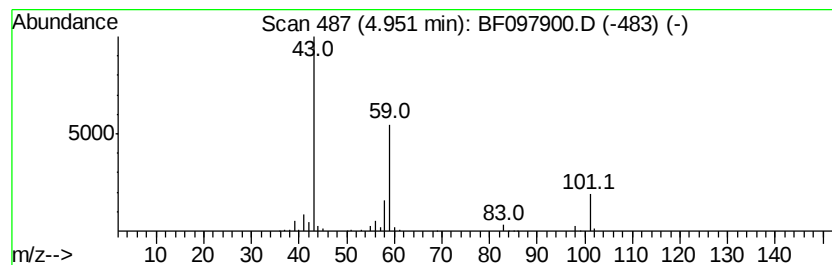
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	11.39 ng	457701	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl ester	141	C7H11NO2	001116-98-9	37
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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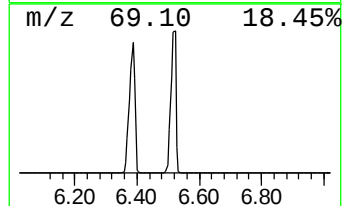
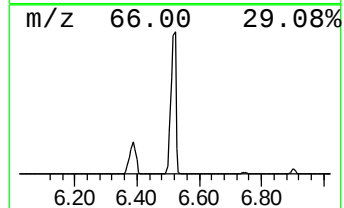
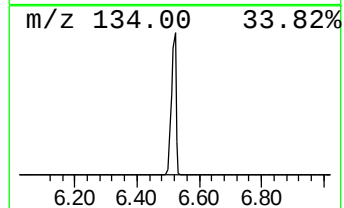
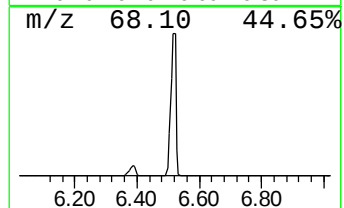
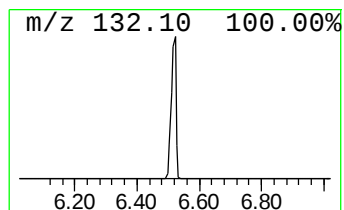
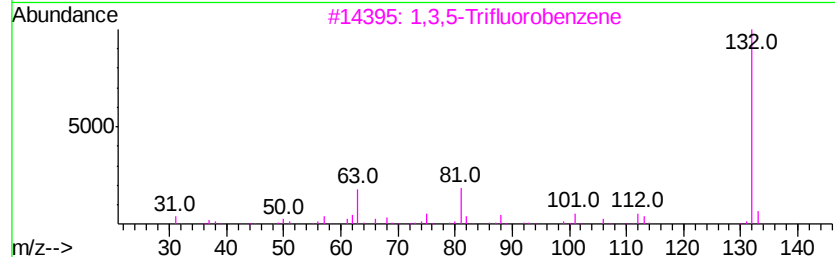
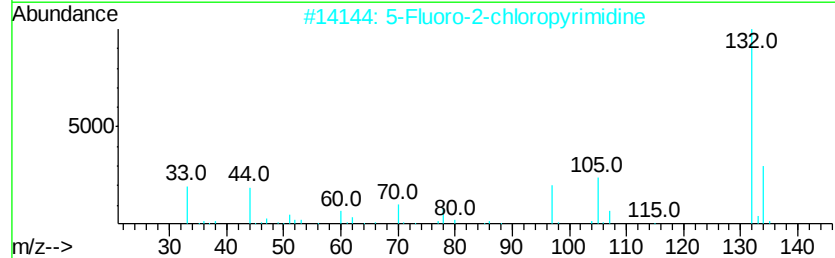
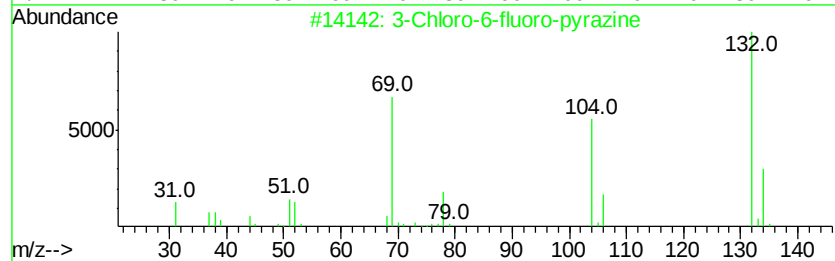
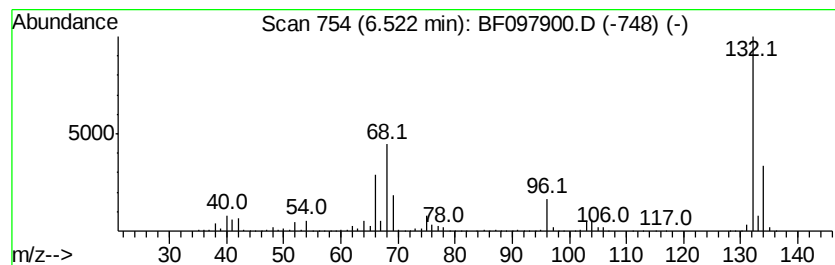
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	93.32 ng	3749600	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4		Xenon	132	Xe	007440-63-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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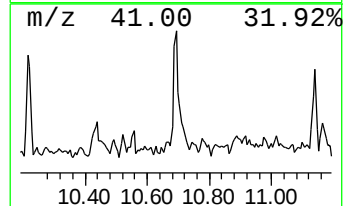
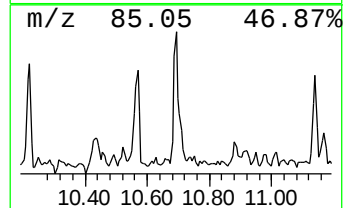
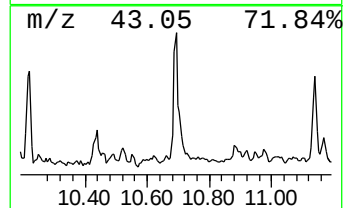
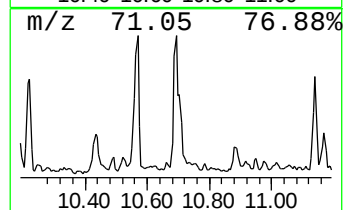
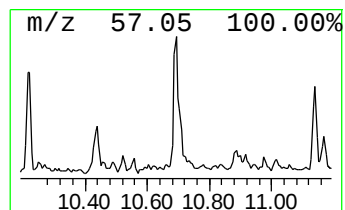
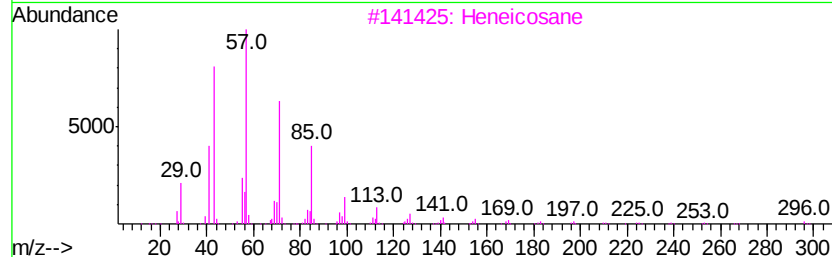
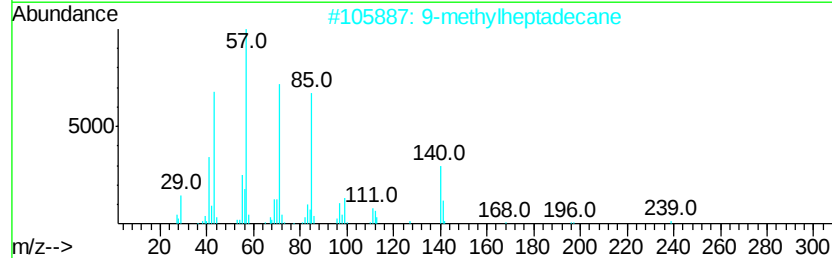
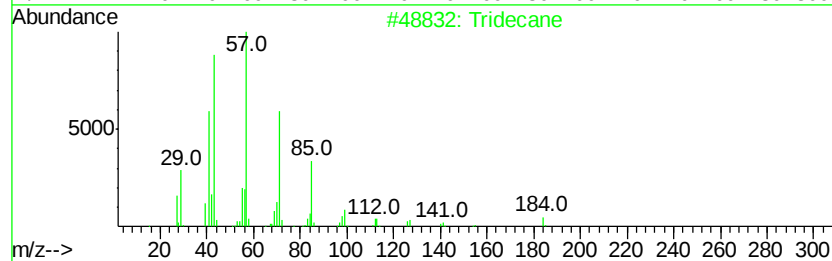
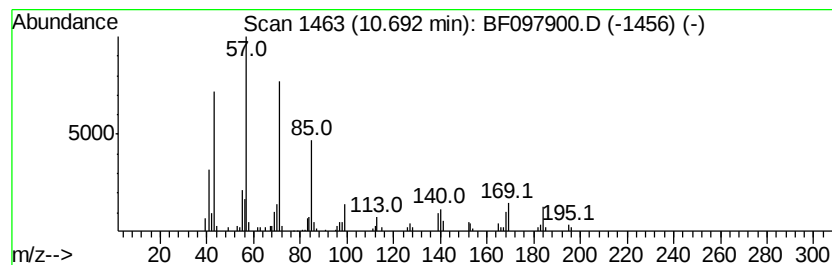
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Tridecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.69	2.61 ng	124650	Phenanthrene-d10		11.26
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	81
2	9-methylheptadecane	254	C18H38	026741-18-4	72
3	Heneicosane	296	C21H44	000629-94-7	68
4	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	64
5	Hexadecane	226	C16H34	000544-76-3	64



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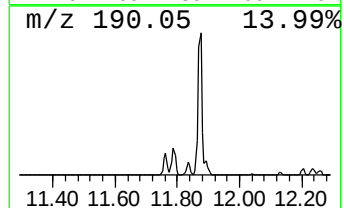
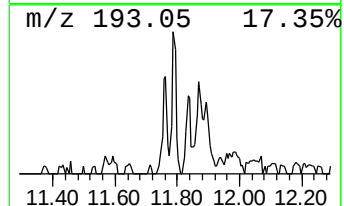
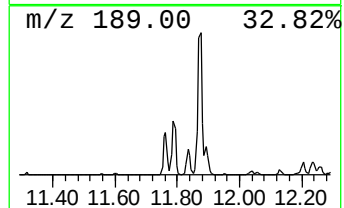
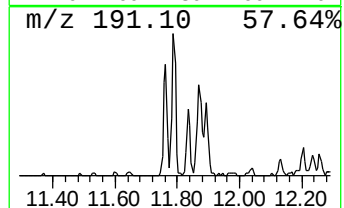
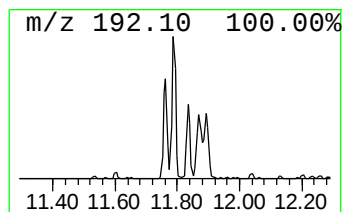
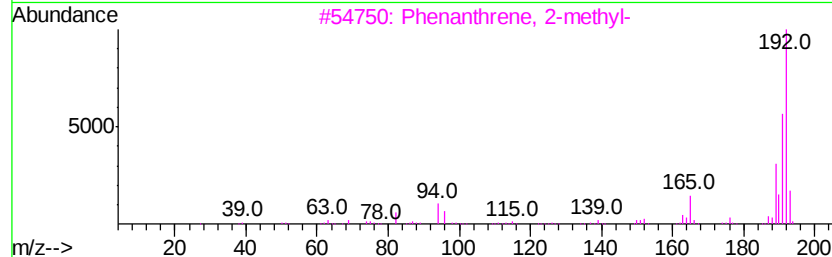
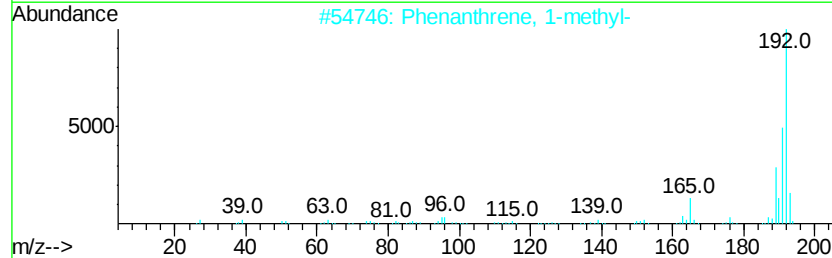
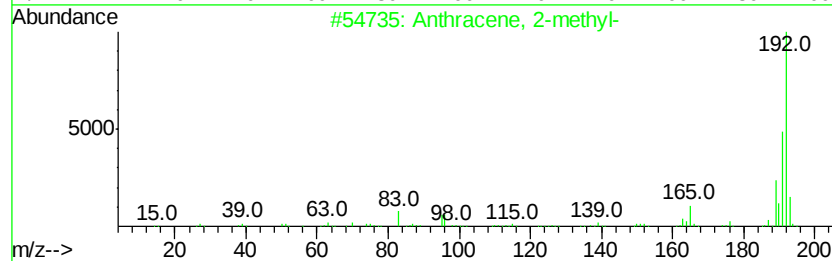
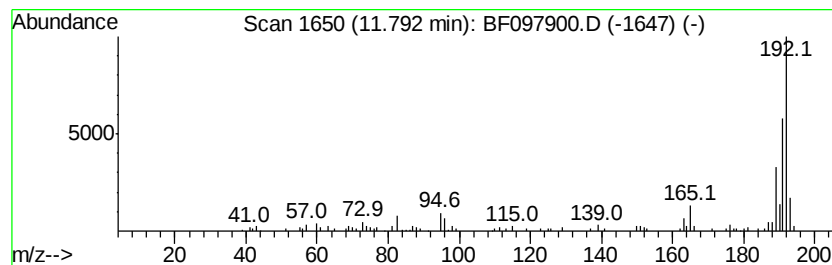
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ClientSampleId :
SASP2P-TP-106-A(0-5)COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.79	2.56 ng	122357	Phenanthrene-d10	11.26	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	96



Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
Data File : BF097900.D
Acq On : 21 Aug 2017 12:18
Operator : SJ/JU
Sample : I4869-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SASP2P-TP-106-A(0-5)COMP

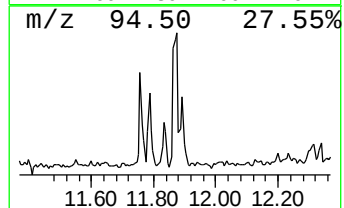
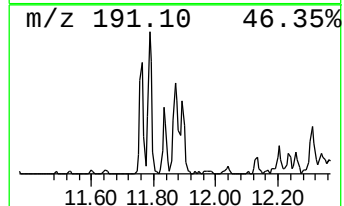
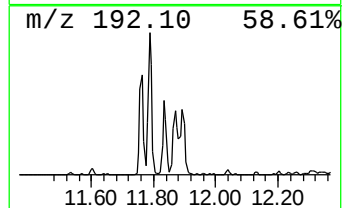
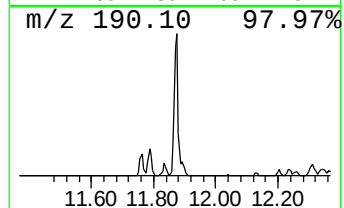
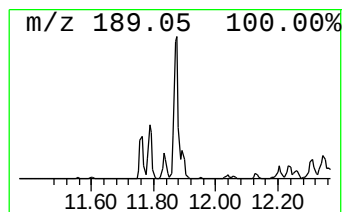
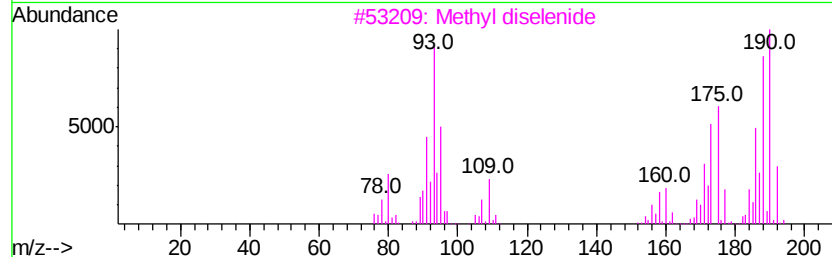
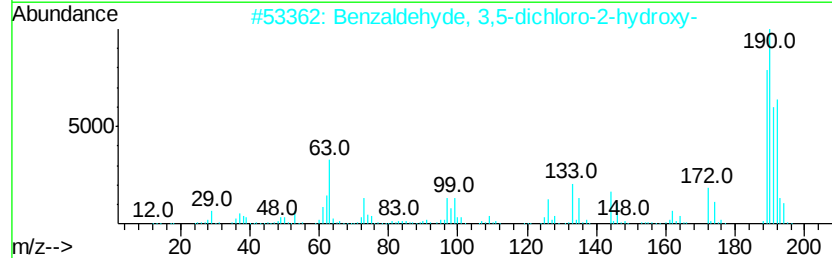
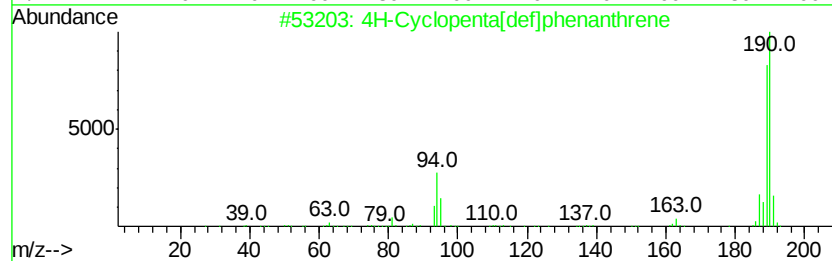
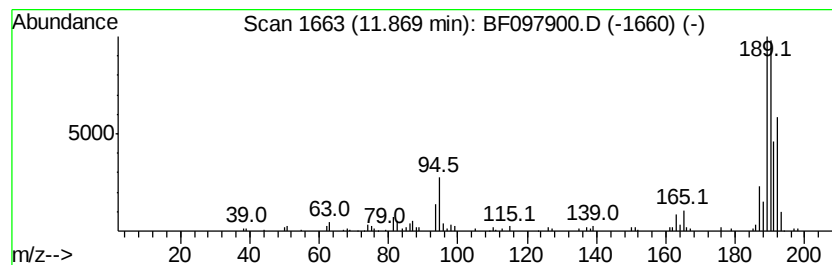
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	3.44 ng	164414	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	47
3		Methyl diselenide	190	C2H6Se2	007101-31-7	38
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	20
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	18



Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
Data File : BF097900.D
Acq On : 21 Aug 2017 12:18
Operator : SJ/JU
Sample : I4869-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SASP2P-TP-106-A(0-5)COMP

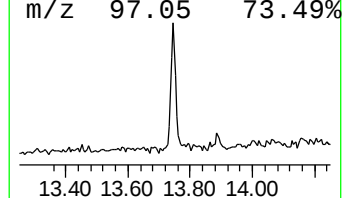
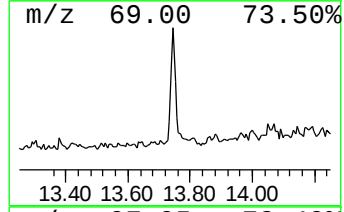
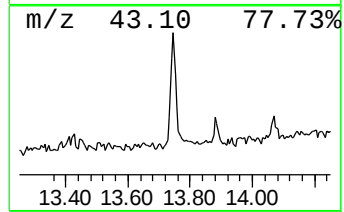
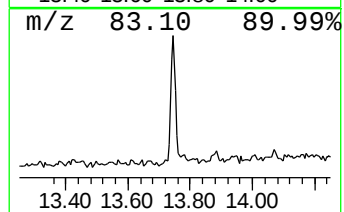
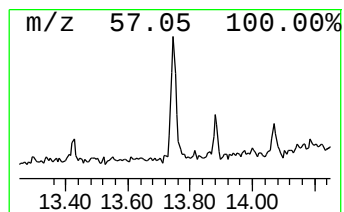
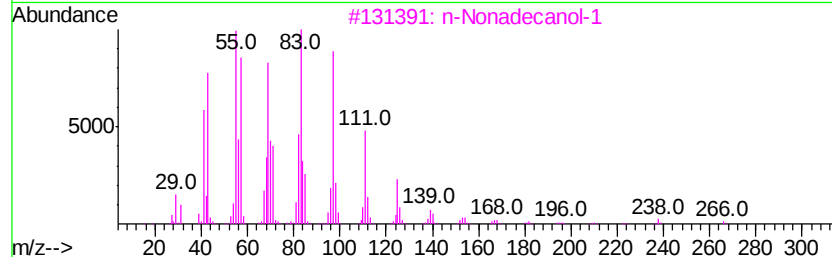
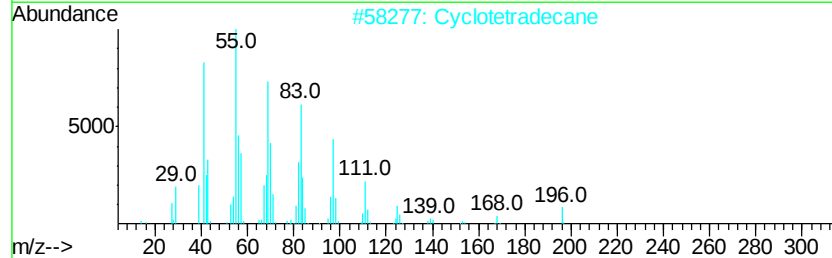
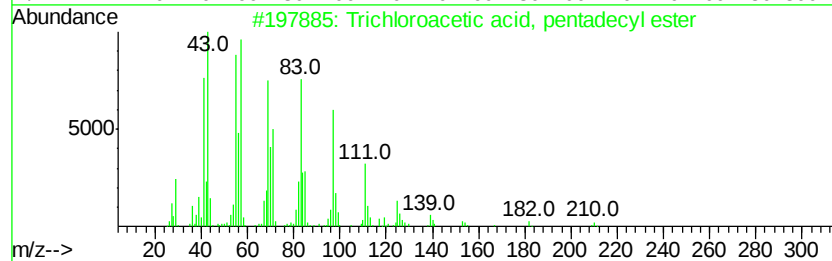
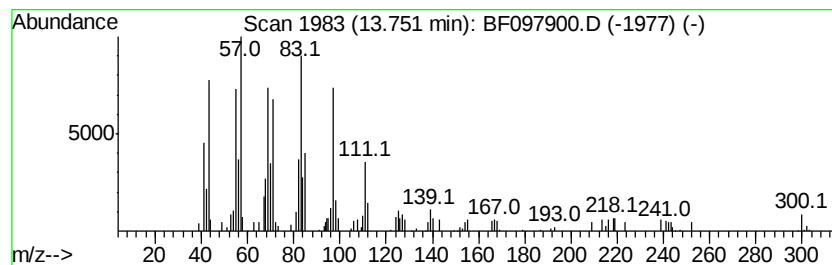
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Trichloroacetic acid, penta... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	3.35 ng	208945	Chrysene-d12	13.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90
2		Cyclotetradecane	196	C14H28	000295-17-0	90
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	87
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	87
5		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	87



Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
Data File : BF097900.D
Acq On : 21 Aug 2017 12:18
Operator : SJ/JU
Sample : I4869-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

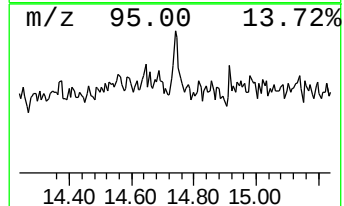
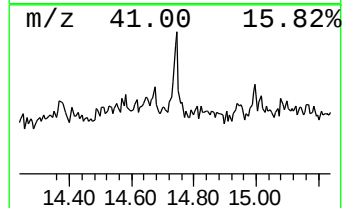
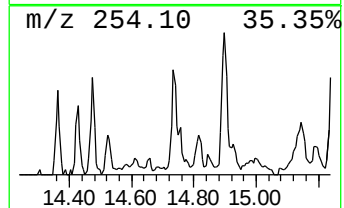
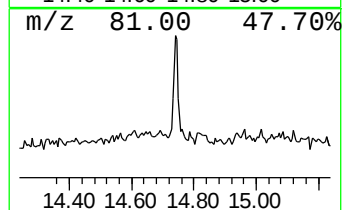
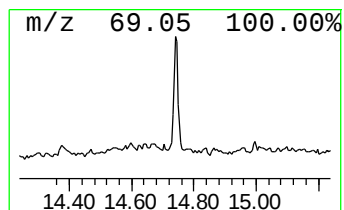
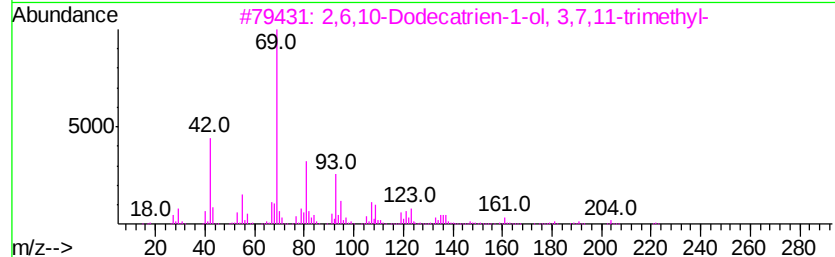
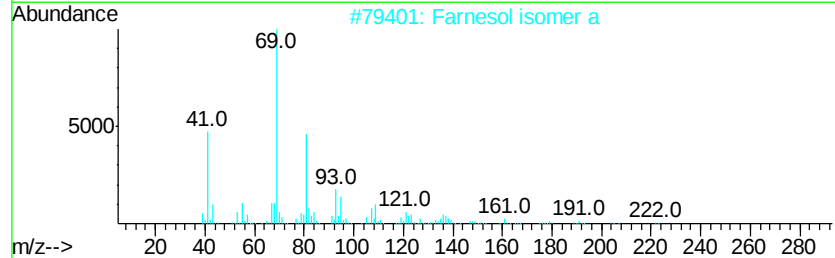
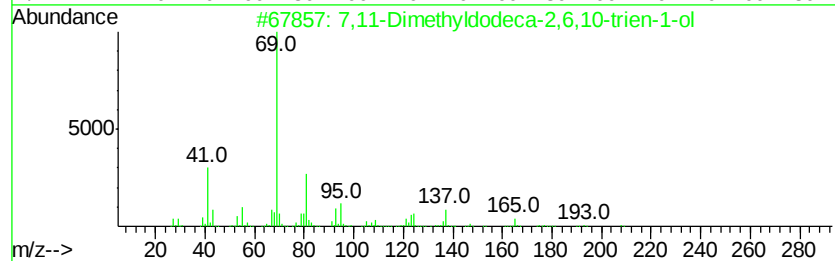
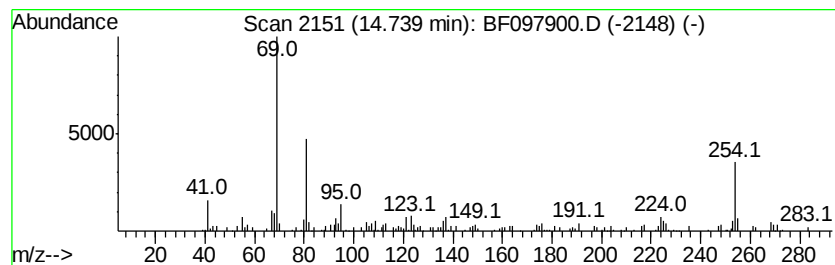
Instrument :
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ClientSampleId :
SASP2P-TP-106-A(0-5)COMP

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7,11-Dimethyldodeca-2,6,10-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.74	2.51 ng	88661	Perylene-d12	15.31	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	58
2	Farnesol isomer a	222	C15H26O	1000108-92-4	58
3	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	52
4	2,6-Octadiene, 2,6-dimethyl-	138	C10H18	002792-39-4	50
5	Trideca-1,3,7,11-tetraene-1,1-di...	268	C18H24N2	1000159-88-9	47



Data Path : U:\HPCHEM1\BNA F\DATA\BF082117\
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Acq On : 21 Aug 2017 12:18
Operator : SJ/JU
Sample : I4869-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

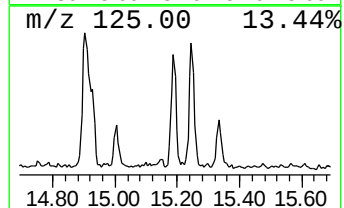
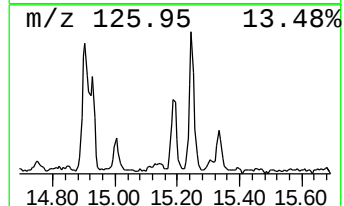
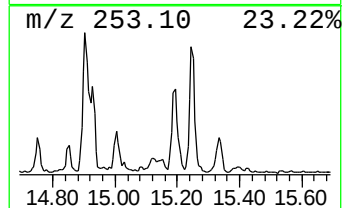
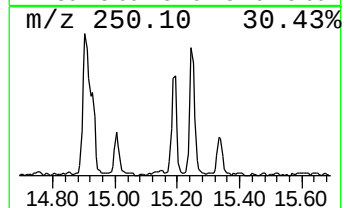
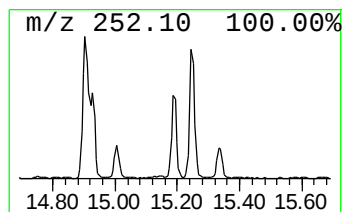
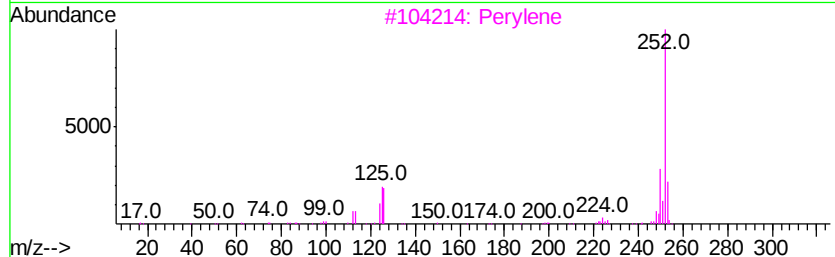
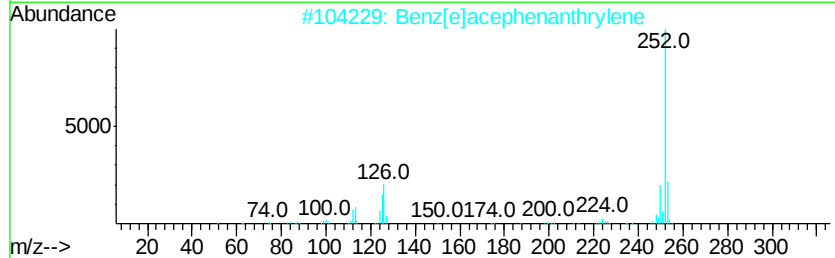
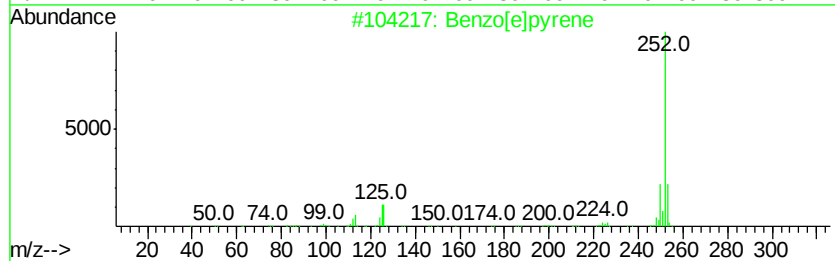
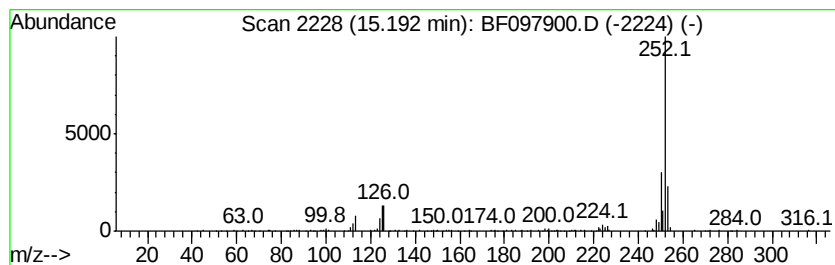
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.19	5.98 ng	210989	Perylene-d12	15.31	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]bpvrene	252	C20H12	000192-97-2	97
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94
3	Pervlene	252	C20H12	000198-55-0	94
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90



Data Path : U:\HPCHEM1\BNA_F\DATA\BF082117\
Data File : BF097900.D
Acq On : 21 Aug 2017 12:18
Operator : SJ/JU
Sample : I4869-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SASP2P-TP-106-A(0-5)COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.13	3.13	2.5	ng	101539	1	6.75	803573	20.0
2-Pentanone, 4-hy...	4.95	11.4	ng	457701	1	6.75	803573	20.0
unknown6.52	6.52	93.3	ng	3749600	1	6.75	803573	20.0
Tridecane	10.69	2.6	ng	124650	4	11.26	956426	20.0
Anthracene, 2-met...	11.79	2.6	ng	122357	4	11.26	956426	20.0
4H-Cyclopenta[def...	11.87	3.4	ng	164414	4	11.26	956426	20.0
Trichloroacetic a...	13.75	3.4	ng	208945	5	13.89	1248940	20.0
7,11-Dimethyldode...	14.74	2.5	ng	88661	6	15.31	705390	20.0
Benzo[e]pyrene	15.19	6.0	ng	210989	6	15.31	705390	20.0