

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000864.D
 Acq On : 18 Oct 2019 03:29
 Operator : JU
 Sample : K5401-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR200036-RT1488

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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP100119.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	5.375	555	559	582	rBV	3489081	3632984	65.14%	8.109%
2	6.393	728	732	750	rBV	3543647	3820214	68.50%	8.527%
3	6.522	750	754	757	rBV	4540182	3965223	71.10%	8.851%
4	6.746	789	792	802	rBV	1198381	1049368	18.82%	2.342%
5	6.905	815	819	822	rBV	4088558	3414028	61.22%	7.620%
6	7.310	884	888	891	rBV	2756428	2413083	43.27%	5.386%
7	8.034	1007	1011	1014	rBV	1514091	1289150	23.12%	2.877%
8	9.116	1191	1195	1198	rBV	6403587	5342945	95.80%	11.926%
9	9.510	1259	1262	1271	rBV	650865	517805	9.28%	1.156%
10	9.799	1307	1311	1314	rBV	1731232	1551828	27.83%	3.464%
11	10.593	1442	1446	1459	rBV	3731766	3343790	59.96%	7.464%
12	11.293	1562	1565	1568	rBV	1943606	1645784	29.51%	3.673%
13	11.316	1568	1569	1574	rVV	260140	194734	3.49%	0.435%
14	11.369	1574	1578	1581	rVB2	101169	120176	2.15%	0.268%
15	11.834	1654	1657	1661	rBV2	101786	109882	1.97%	0.245%
16	12.522	1770	1774	1777	rBV	709882	598271	10.73%	1.335%
17	12.751	1808	1813	1816	rBV	560023	538464	9.66%	1.202%
18	12.904	1834	1839	1842	rBV	6200136	5576930	100.00%	12.448%
19	13.087	1868	1870	1873	rVB	82960	76204	1.37%	0.170%
20	13.822	1992	1995	2005	rVB3	398894	520871	9.34%	1.163%
21	13.969	2015	2020	2023	rBV	1731389	1821218	32.66%	4.065%
22	13.998	2023	2025	2028	rVB	186375	149414	2.68%	0.333%
23	14.075	2036	2038	2042	rVB	79677	77270	1.39%	0.172%
24	14.145	2048	2050	2056	rVB2	95890	96427	1.73%	0.215%
25	14.457	2100	2103	2105	rBV3	120909	133841	2.40%	0.299%
26	14.616	2127	2130	2134	rBV	115028	144768	2.60%	0.323%
27	14.763	2153	2155	2158	rVB	131433	117239	2.10%	0.262%
28	15.022	2196	2199	2207	rVB	197106	348969	6.26%	0.779%
29	15.104	2210	2213	2216	rBV	134724	134033	2.40%	0.299%
30	15.322	2248	2250	2257	rVB	104756	127014	2.28%	0.284%
31	15.387	2258	2261	2267	rVB	135402	161146	2.89%	0.360%
32	15.451	2267	2272	2276	rBV	1364675	1629358	29.22%	3.637%
33	15.928	2349	2353	2359	rVB	100064	139374	2.50%	0.311%

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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-BP100119.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

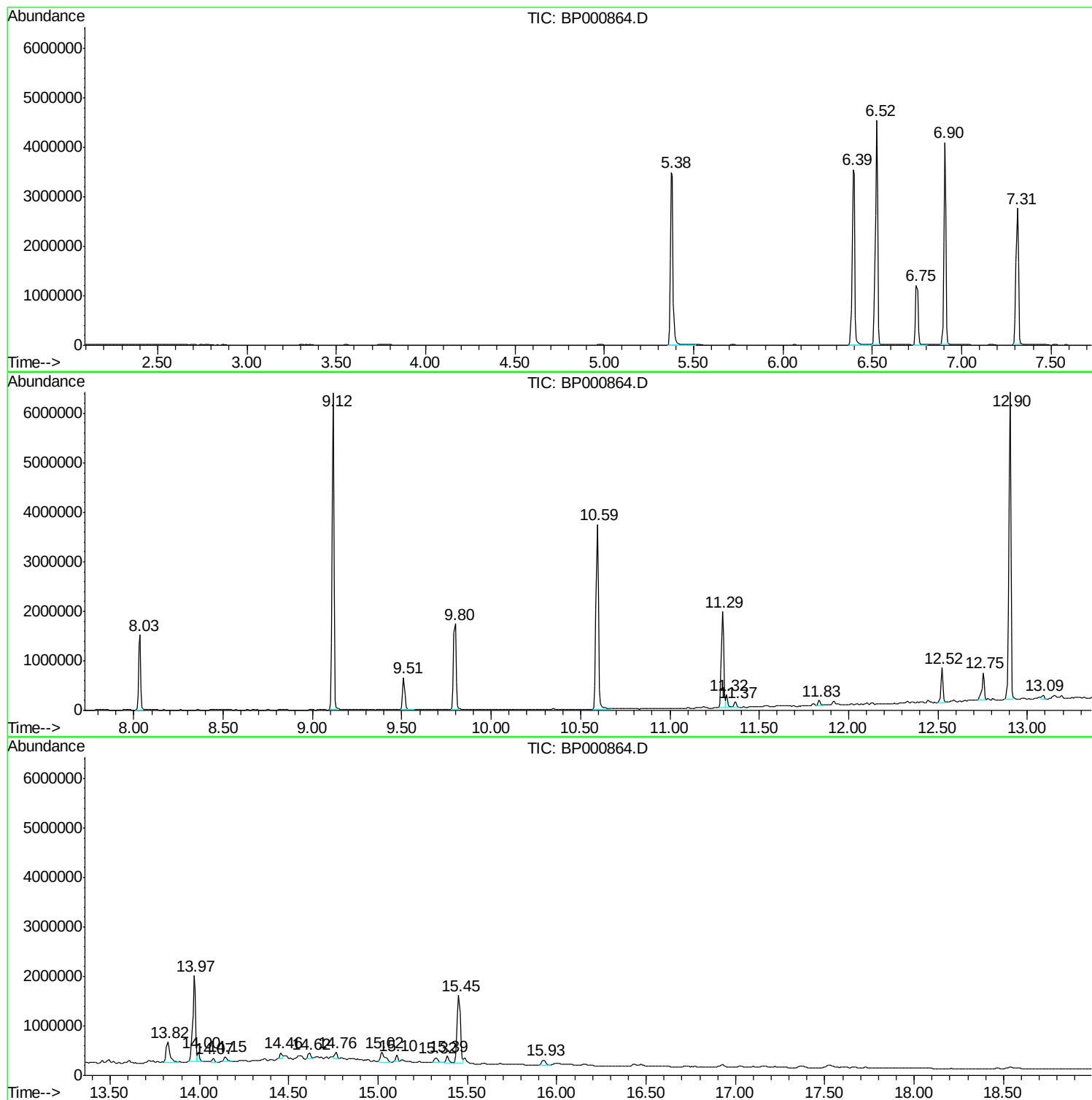
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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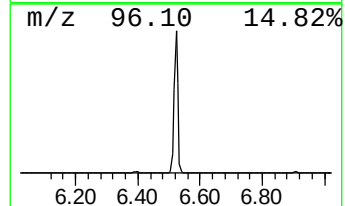
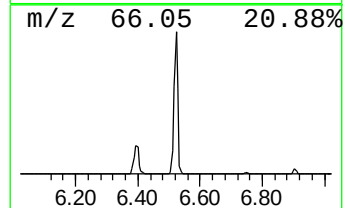
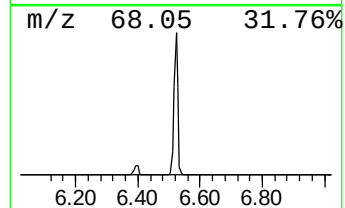
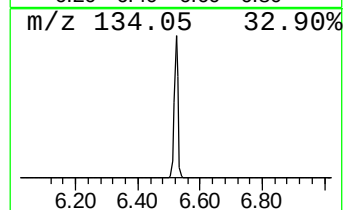
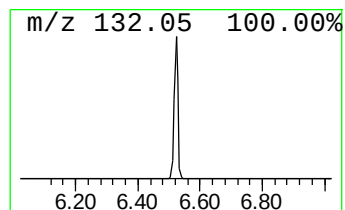
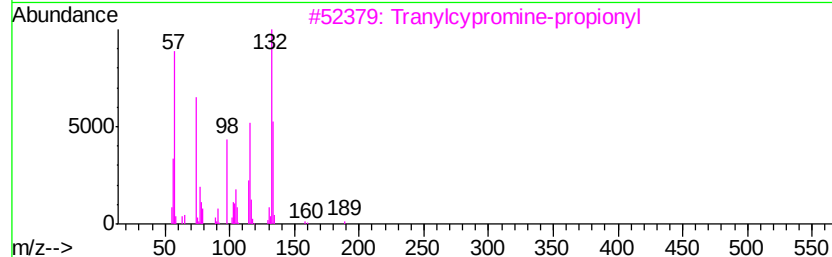
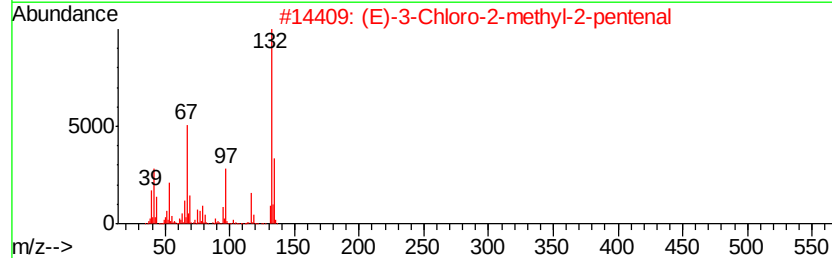
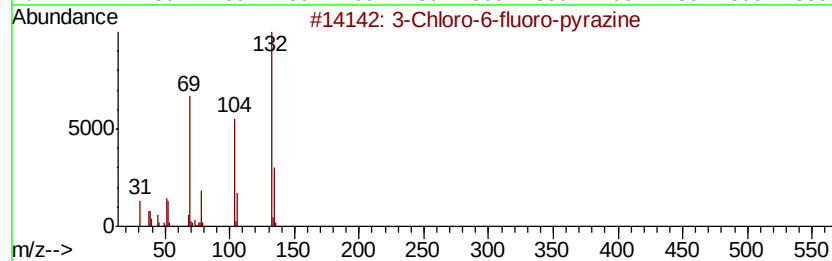
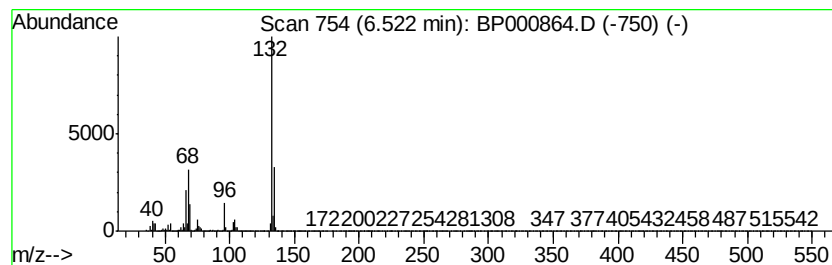
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Peak Number 1 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	75.57 ng	3965220	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	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	16
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	12



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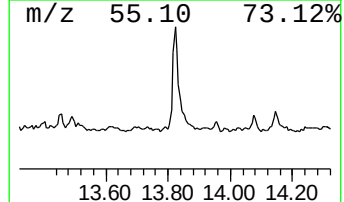
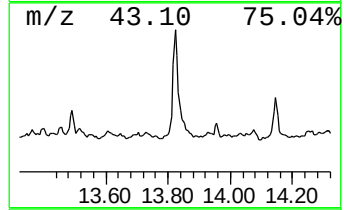
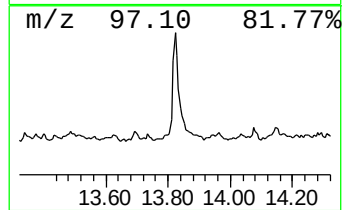
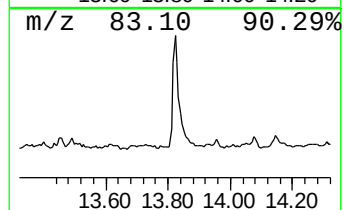
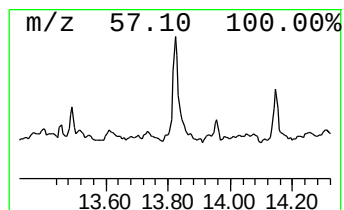
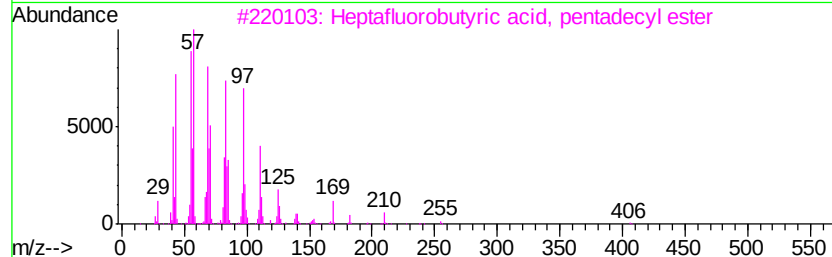
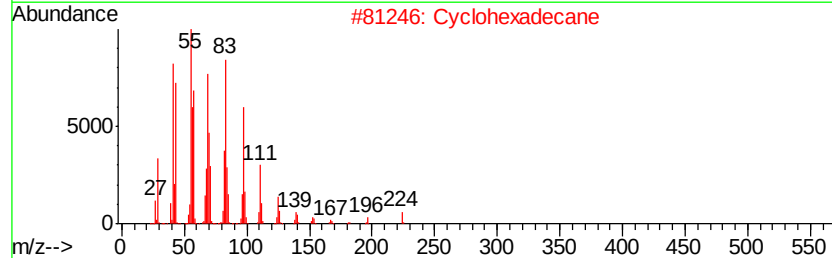
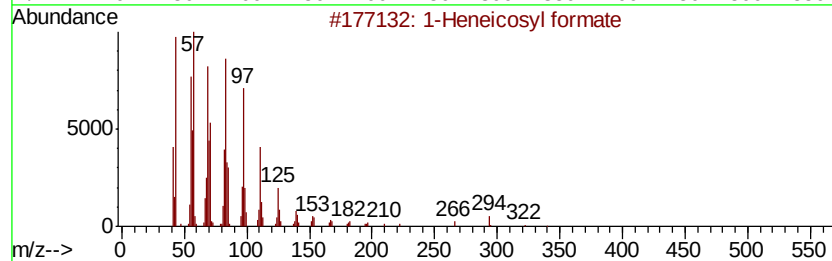
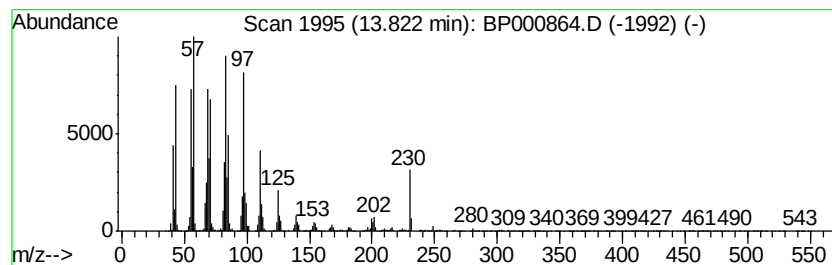
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Peak Number 3 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.82	5.72 ng	520871	Chrysene-d12		13.97
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2	Cyclohexadecane	224	C16H32	000295-65-8	93
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
4	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown6.52	6.52	75.6	ng	3965220	1	6.75	1049370	20.0
1-Heneicosyl formate	13.82	5.7	ng	520871	5	13.97	1821220	20.0