

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032822\
 Data File : BF127555.D
 Acq On : 28 Mar 2022 21:28
 Operator : CG\JU
 Sample : N2108-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R206

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-BF032122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127555.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.440	186	196	213	rBV	254201	801464	34.49%	4.711%
2	5.081	472	475	486	rBV	20207	33567	1.44%	0.197%
3	5.463	536	540	560	rBV	2289895	2302775	99.11%	13.536%
4	6.475	708	712	732	rBV	2414718	2323517	100.00%	13.658%
5	6.834	769	773	781	rBV	615343	577061	24.84%	3.392%
6	7.004	800	802	810	rVB	39099	37376	1.61%	0.220%
7	7.398	865	869	881	rBV	1606421	1581460	68.06%	9.296%
8	8.116	987	991	994	rBV	800834	761727	32.78%	4.478%
9	8.140	994	995	1001	rVB	74509	46797	2.01%	0.275%
10	8.834	1110	1113	1120	rVB	39018	38039	1.64%	0.224%
11	8.928	1126	1129	1135	rVB	27828	23784	1.02%	0.140%
12	9.187	1169	1173	1177	rBV	2308218	1889630	81.33%	11.108%
13	9.869	1284	1289	1293	rBV	1028119	835070	35.94%	4.909%
14	9.904	1293	1295	1301	rVB	59751	49354	2.12%	0.290%
15	10.081	1320	1325	1328	rBV	38105	34137	1.47%	0.201%
16	10.422	1378	1383	1387	rBV	48944	42454	1.83%	0.250%
17	10.663	1420	1424	1434	rBV	1565073	1352149	58.19%	7.948%
18	11.363	1539	1543	1545	rBV	897444	763964	32.88%	4.491%
19	11.386	1545	1547	1551	rVB	174486	140610	6.05%	0.827%
20	11.610	1581	1585	1591	rBV2	22231	24692	1.06%	0.145%
21	11.880	1626	1631	1632	rBV2	32699	41603	1.79%	0.245%
22	11.980	1645	1648	1650	rBV2	26491	24485	1.05%	0.144%
23	12.580	1747	1750	1754	rBV	174199	146322	6.30%	0.860%
24	12.810	1783	1789	1794	rBV	116087	128844	5.55%	0.757%
25	12.945	1808	1812	1815	rBV	1427437	1329858	57.23%	7.817%
26	13.769	1945	1952	1954	rBV2	16041	23376	1.01%	0.137%
27	13.845	1958	1965	1978	rVV5	66774	224934	9.68%	1.322%
28	14.010	1989	1993	1997	rBV	634885	615476	26.49%	3.618%
29	14.039	1997	1998	2004	rVB	48652	36078	1.55%	0.212%
30	14.386	2054	2057	2059	rBV4	30348	28913	1.24%	0.170%
31	15.069	2170	2173	2174	rBV	33396	40127	1.73%	0.236%
32	15.374	2222	2225	2231	rBV	21267	31297	1.35%	0.184%
33	15.439	2233	2236	2239	rVV	27941	42347	1.82%	0.249%
34	15.486	2241	2244	2256	rVB	466343	603775	25.99%	3.549%
35	15.892	2309	2313	2319	rBV2	25447	34780	1.50%	0.204%

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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-BF032122.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

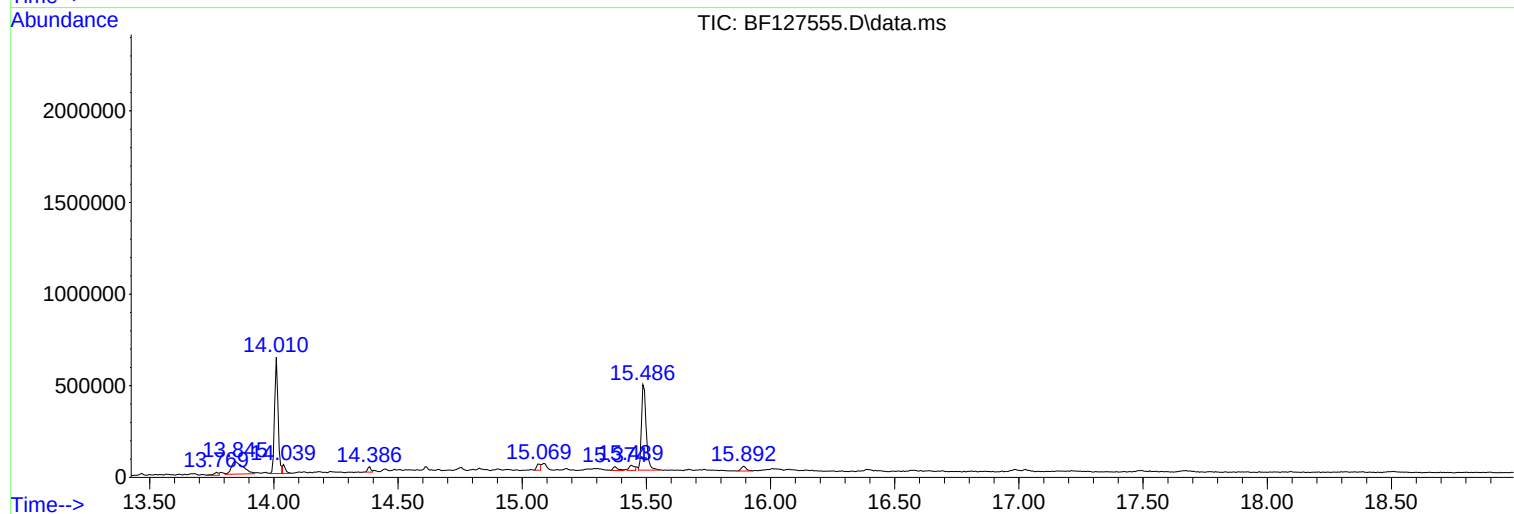
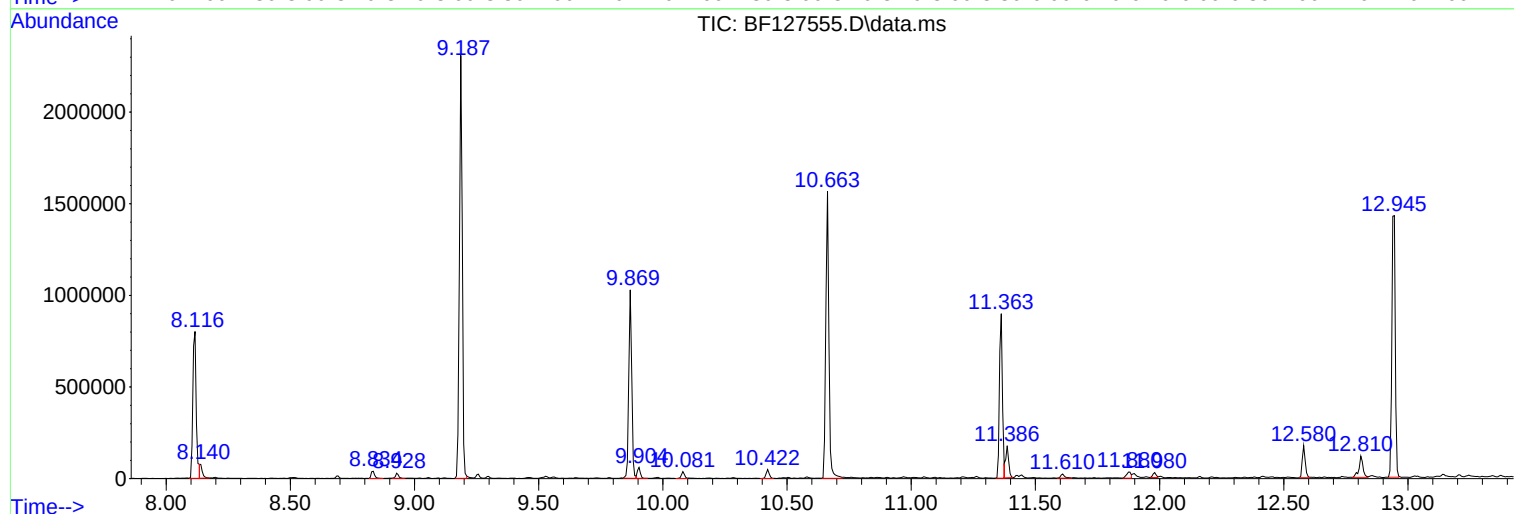
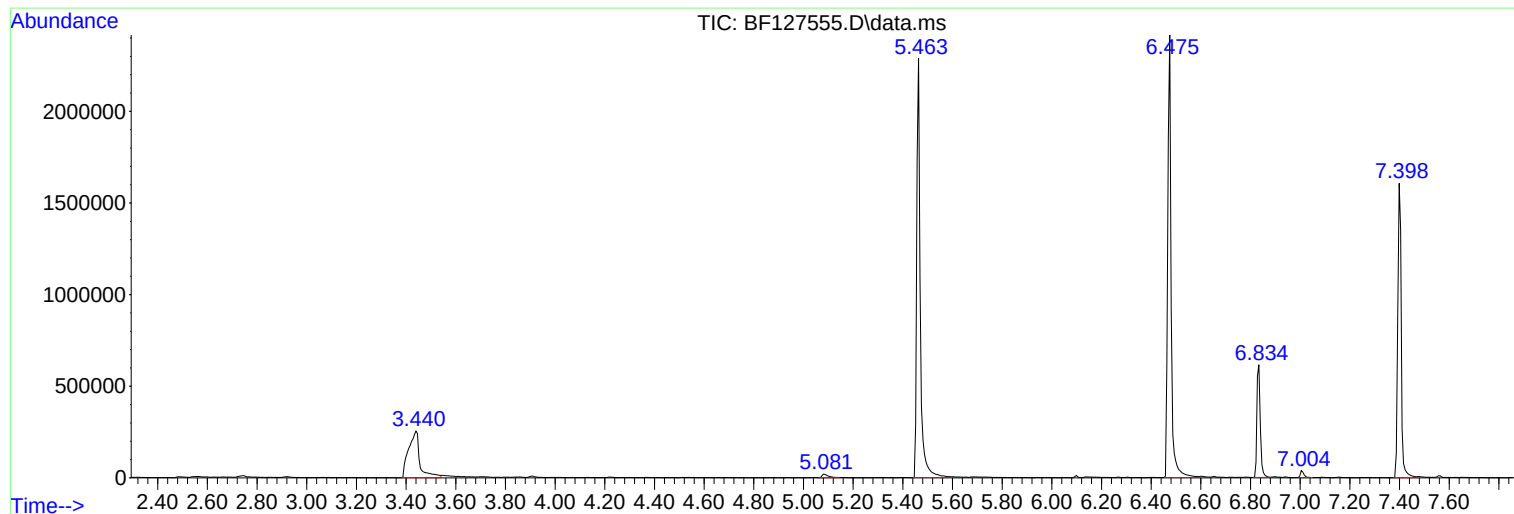
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF032122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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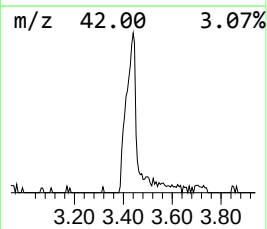
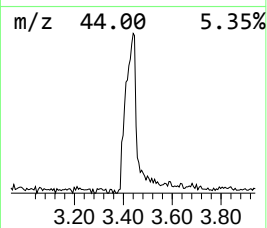
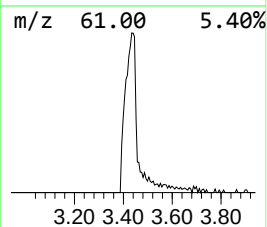
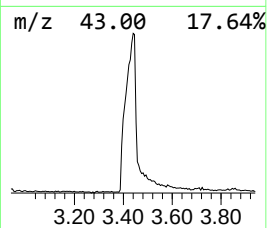
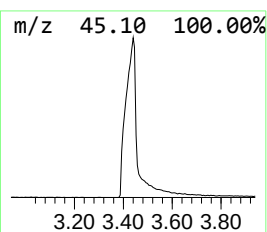
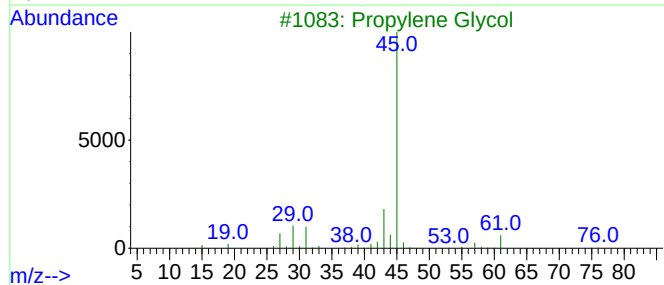
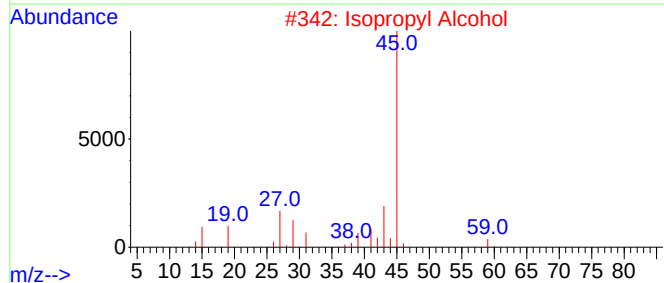
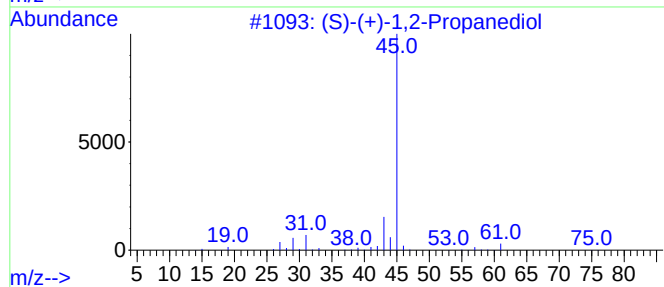
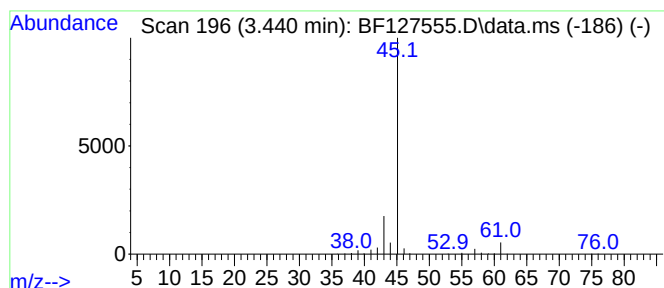
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (S)-(+)-1,2-Propanediol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.440	27.78 ng	801464	1,4-Dichlorobenzene-d4	6.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-1,2-Propanediol			76	C3H8O2	004254-15-3	64
2	Isopropyl Alcohol			60	C3H8O	000067-63-0	56
3	Propylene Glycol			76	C3H8O2	000057-55-6	47
4	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	43
5	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	9



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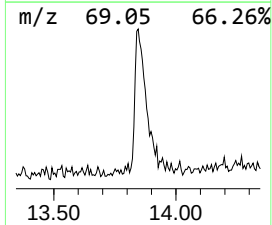
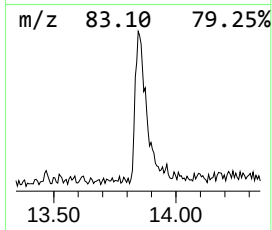
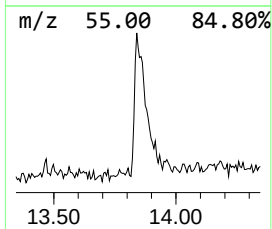
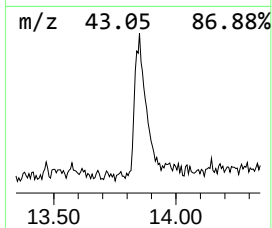
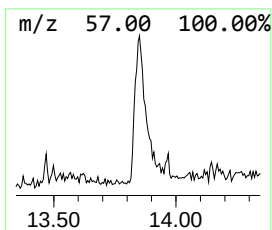
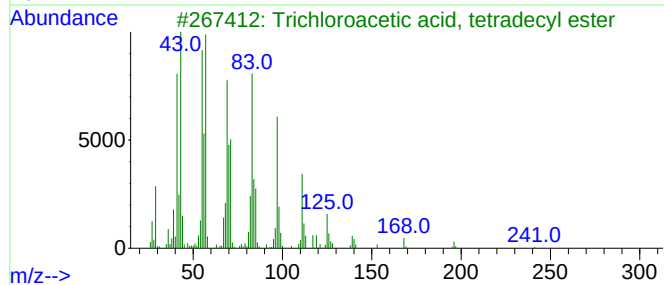
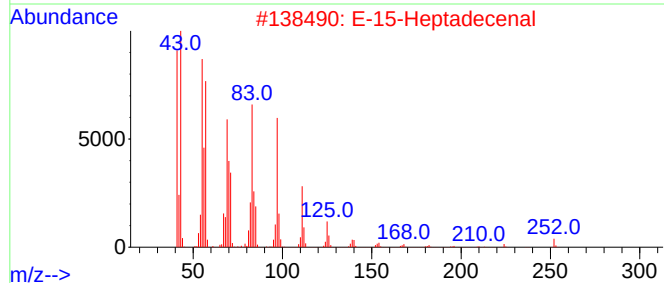
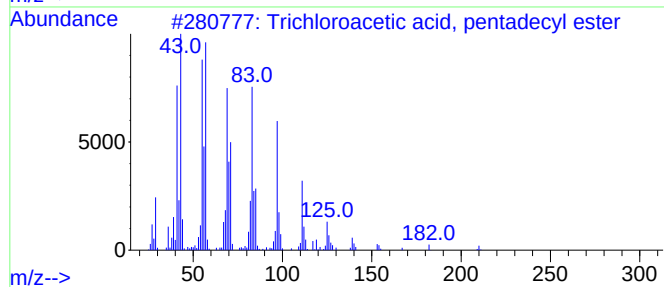
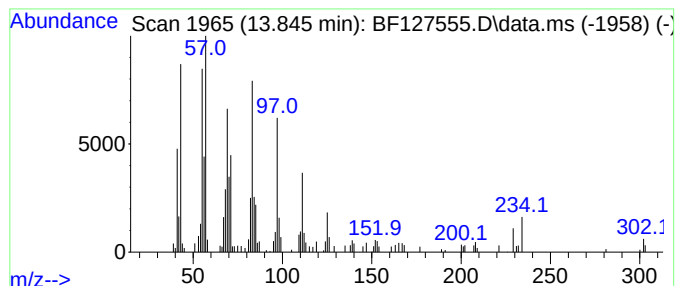
Instrument :
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Peak Number 2 Trichloroacetic acid, penta... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.845	7.31 ng	224934	Chrysene-d12		14.010	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Trichloroacetic acid, pentadecyl...		372	C17H31Cl3O2	074339-53-0	90
2	E-15-Heptadecenal		252	C17H32O	1000130-97-9	90
3	Trichloroacetic acid, tetradecyl...		358	C16H29Cl3O2	074339-52-9	90
4	1-Hexacosene		364	C26H52	018835-33-1	90
5	Bromoacetic acid, octadecyl ester		390	C20H39BrO2	018992-03-5	87



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(S)-(+)-1,2-Pro...	3.440	27.8	ng	801464	1	6.834	577061	20.0
Trichloroacetic...	13.845	7.3	ng	224934	5	14.010	615476	20.0