

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071521\
 Data File : BP006235.D
 Acq On : 15 Jul 2021 12:59
 Operator : CG/JU
 Sample : PB137674BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB137674BL

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_P\Methods\8270E-BP070721.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006235.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.422	391	397	423	rBV	2343405	3962605	53.35%	11.157%
2	7.034	665	671	702	rBV	1926594	3850289	51.84%	10.841%
3	7.840	802	808	822	rBV	455218	783161	10.54%	2.205%
4	9.016	1000	1008	1032	rBV	1124507	2361890	31.80%	6.650%
5	10.652	1278	1286	1305	rBV	390061	955208	12.86%	2.689%
6	13.104	1690	1703	1730	rBV	3255282	5337159	71.85%	15.027%
7	14.481	1930	1937	1952	rBV	683985	1252669	16.86%	3.527%
8	15.981	2184	2192	2221	rBV	2039303	3793898	51.08%	10.682%
9	17.234	2398	2405	2425	rBV	720855	1415018	19.05%	3.984%
10	19.481	2783	2787	2790	rBV	96863	111905	1.51%	0.315%
11	19.781	2832	2838	2850	rBV	6160681	7427793	100.00%	20.914%
12	20.975	3036	3041	3045	rBV2	129637	236164	3.18%	0.665%
13	21.028	3047	3050	3055	rVB	180142	235546	3.17%	0.663%
14	21.086	3056	3060	3068	rBV	444517	715022	9.63%	2.013%
15	21.310	3094	3098	3109	rBV	1017670	1443587	19.43%	4.065%
16	23.692	3497	3503	3514	rVB2	731863	1634496	22.01%	4.602%

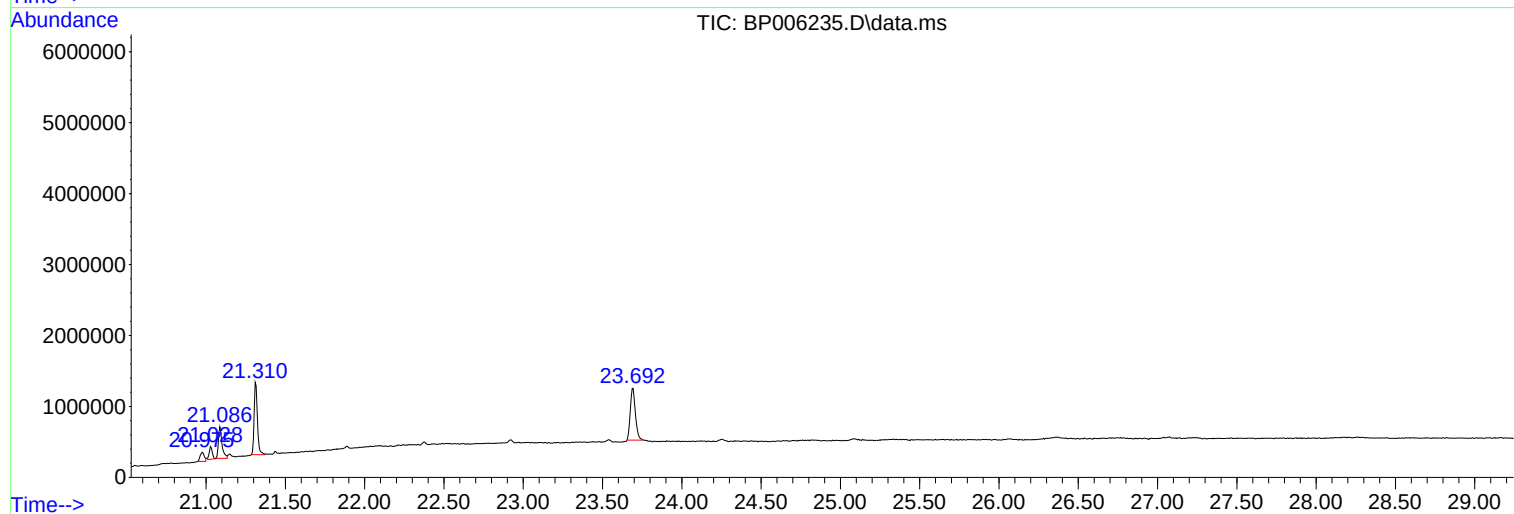
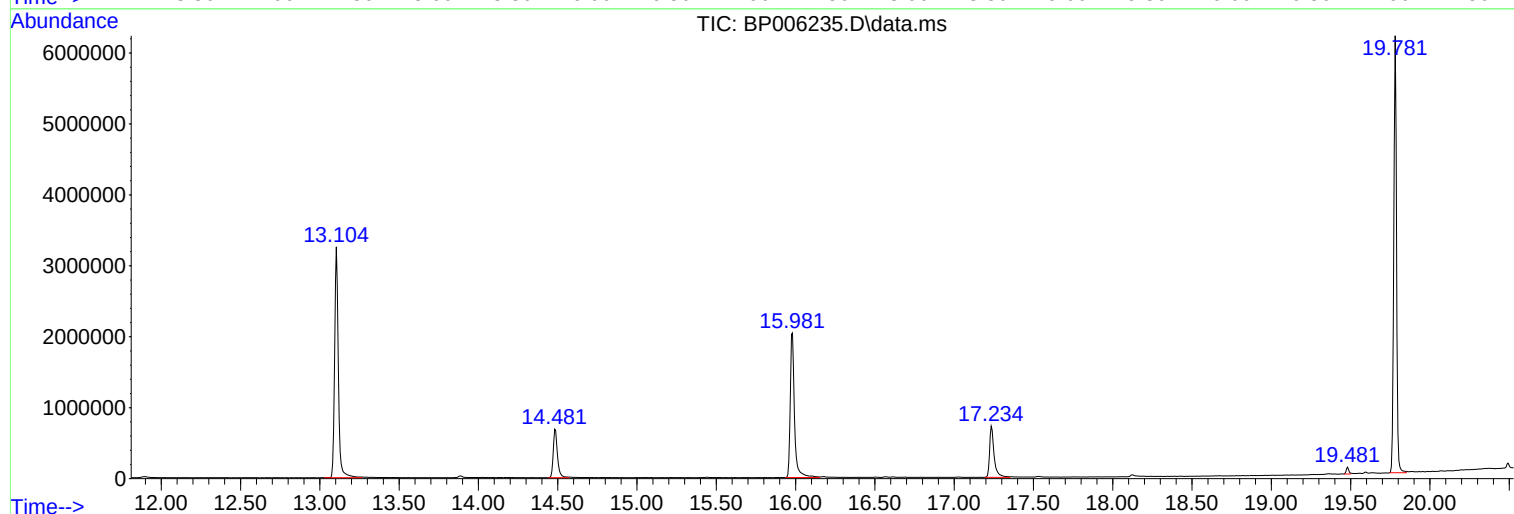
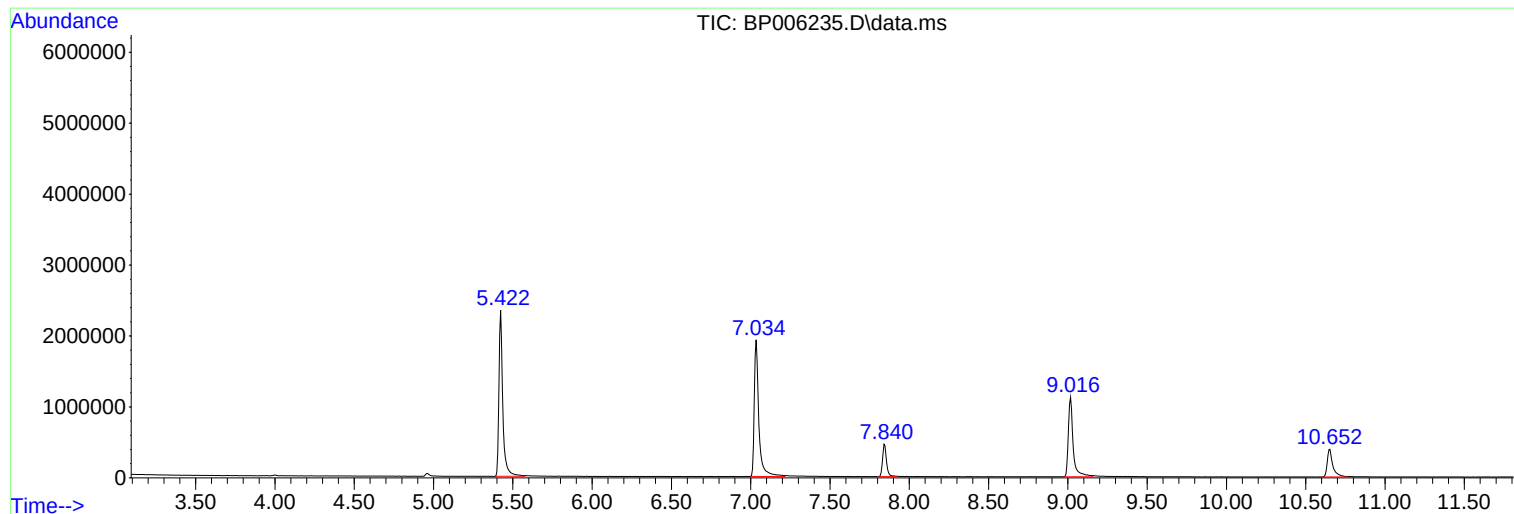
Sum of corrected areas: 35516410

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.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 1 Oxybis(propane-1,2-diyl) di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.975	3.27 ng	236164	Chrysene-d12	21.310

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	53