

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007646.D
 Acq On : 25 Oct 2021 17:41
 Operator : CG/JU
 Sample : PB140190BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK190

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % 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\SFAM-EPA-BP102521.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007646.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	10	14	20	rBV	43008	59018	1.38%	0.143%
2	3.787	81	85	101	rBV	551125	869479	20.28%	2.109%
3	4.122	138	142	146	rVB	9846	12745	0.30%	0.031%
4	7.105	641	649	661	rBV	632211	986077	23.00%	2.392%
5	7.269	670	677	685	rBV	487265	773507	18.04%	1.876%
6	7.469	704	711	721	rBV	638530	1011395	23.59%	2.453%
7	7.940	784	791	800	rVB	533588	830500	19.37%	2.015%
8	8.646	902	911	925	rBV	748597	1226923	28.62%	2.976%
9	9.093	981	987	1002	rVB	552592	929608	21.68%	2.255%
10	9.822	1103	1111	1121	rBV	461962	761040	17.75%	1.846%
11	10.363	1196	1203	1212	rBV	755367	1237246	28.86%	3.001%
12	10.745	1260	1268	1276	rBV	685106	1167681	27.24%	2.832%
13	10.875	1282	1290	1304	rBV	745911	1270657	29.64%	3.082%
14	12.334	1529	1538	1545	rBV2	15206	28891	0.67%	0.070%
15	12.551	1569	1575	1579	rBV9	2094	4759	0.11%	0.012%
16	13.545	1740	1744	1749	rBV8	3551	5772	0.13%	0.014%
17	13.986	1812	1819	1830	rBV	1423318	1929915	45.02%	4.681%
18	14.269	1859	1867	1880	rBV	1497190	2233217	52.09%	5.417%
19	14.575	1911	1919	1933	rBV	1086121	1654252	38.59%	4.013%
20	14.769	1945	1952	1973	rBV	637994	937227	21.86%	2.273%
21	15.569	2081	2088	2098	rBV	1971726	2842656	66.31%	6.895%
22	15.686	2102	2108	2122	rVV	506487	741518	17.30%	1.799%
23	15.810	2125	2129	2135	rVB3	14015	21094	0.49%	0.051%
24	16.139	2180	2185	2189	rVB8	2542	5125	0.12%	0.012%
25	16.263	2200	2206	2207	rBV5	2863	4943	0.12%	0.012%
26	17.016	2330	2334	2337	rVB6	3464	4498	0.10%	0.011%
27	17.116	2346	2351	2354	rBV4	8174	12404	0.29%	0.030%
28	17.327	2380	2387	2397	rBV	1361435	2048378	47.78%	4.969%
29	17.427	2398	2404	2418	rVV2	2135856	3177197	74.11%	7.707%
30	19.245	2708	2713	2717	rBV4	33079	50867	1.19%	0.123%
31	19.310	2720	2724	2727	rBV2	31783	41390	0.97%	0.100%
32	19.663	2778	2784	2793	rBV	3009310	3930971	91.70%	9.535%
33	21.098	3024	3028	3032	rBV	80607	117827	2.75%	0.286%
34	21.151	3033	3037	3041	rBV	131495	189993	4.43%	0.461%
35	21.198	3041	3045	3052	rVV	427855	539682	12.59%	1.309%
36	21.416	3077	3082	3088	rBV	1869997	2564926	59.83%	6.222%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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\SFAM-EPA-BP102521.M
Title : SVOA CALIBRATION

37	23.704	3464	3471	3480	rBV2	2103575	4286930	100.00%	10.399%
38	23.862	3492	3498	3508	rVB2	1324082	2715618	63.35%	6.587%

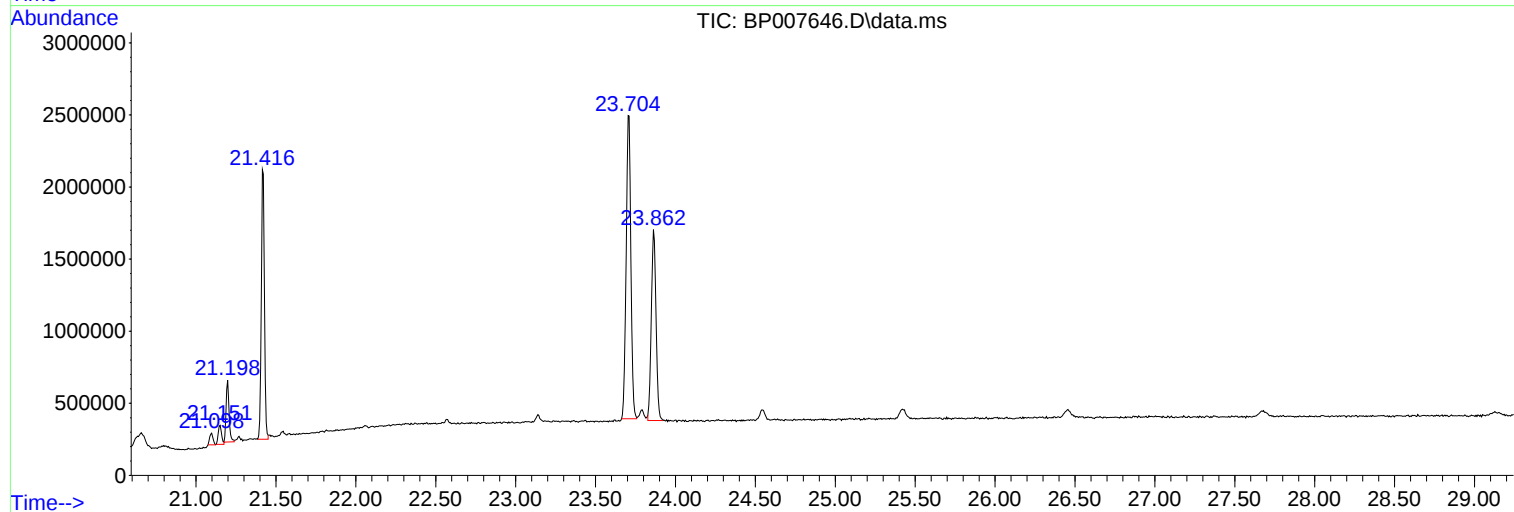
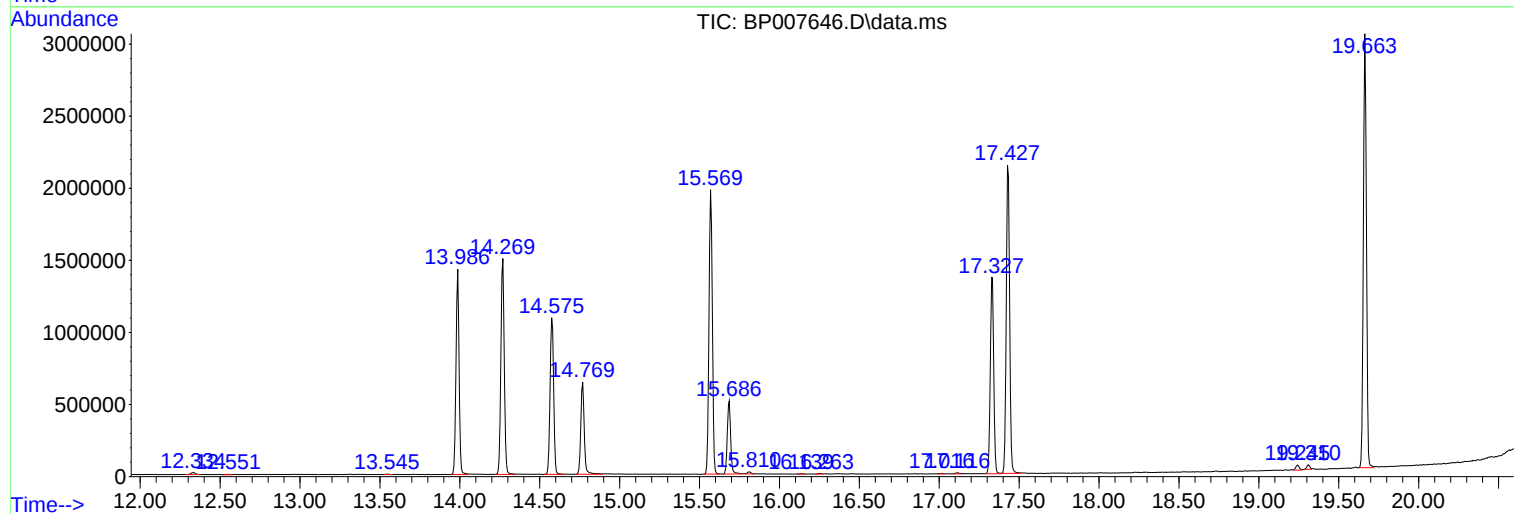
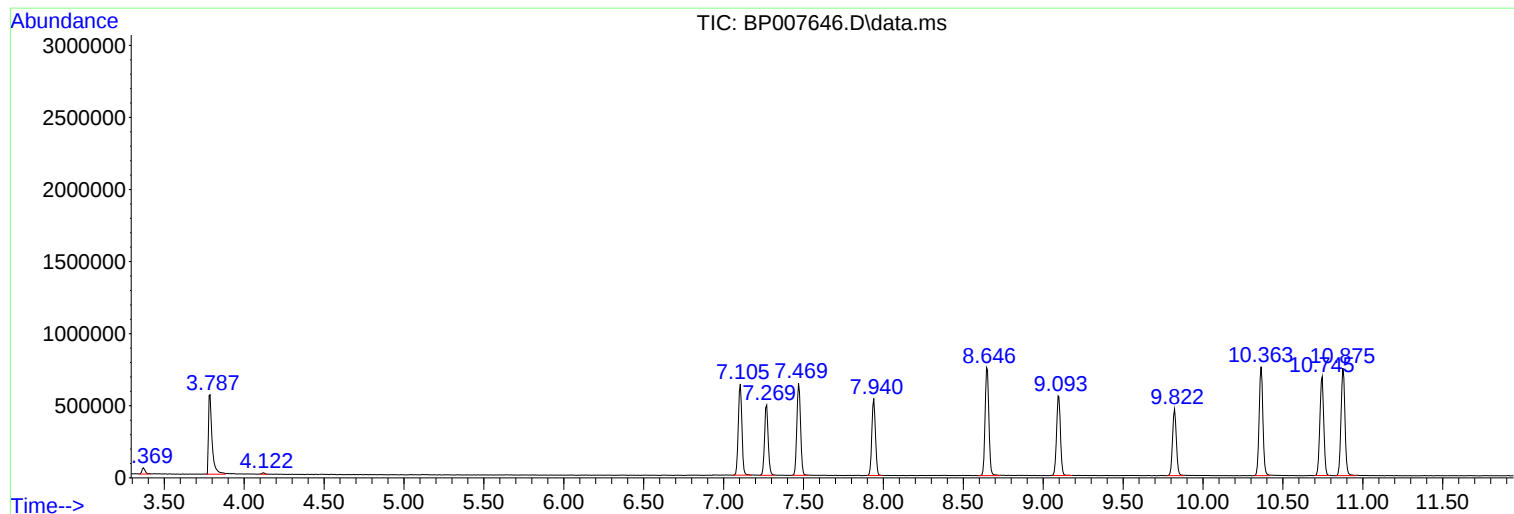
Sum of corrected areas: 41225926

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION

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



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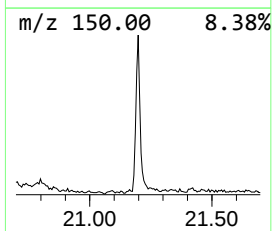
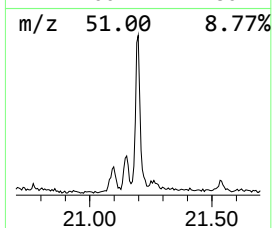
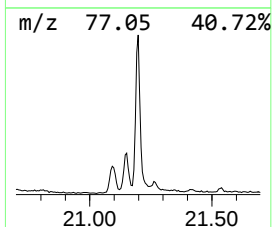
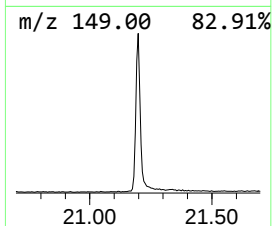
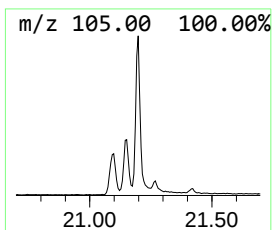
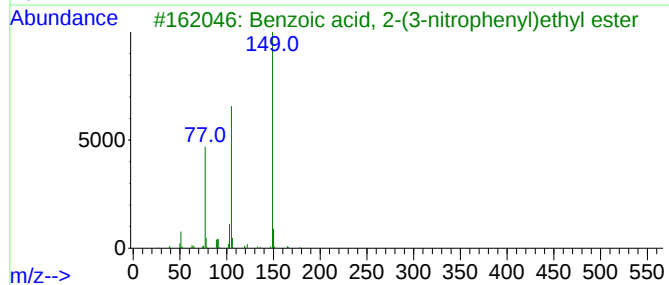
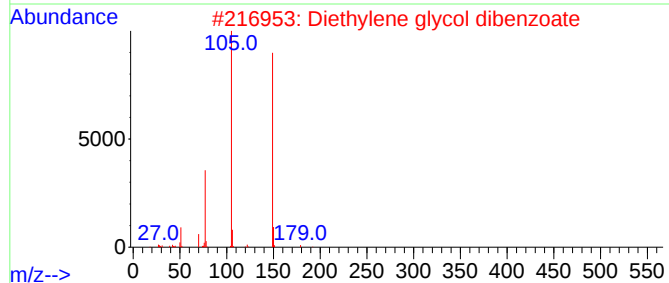
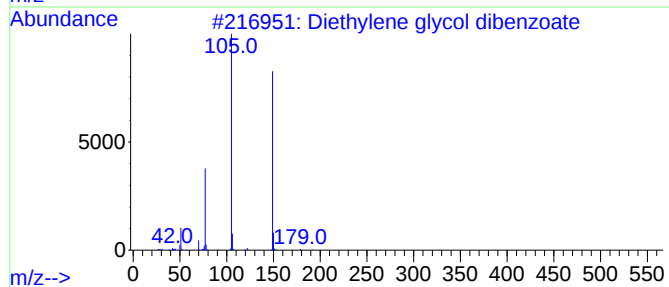
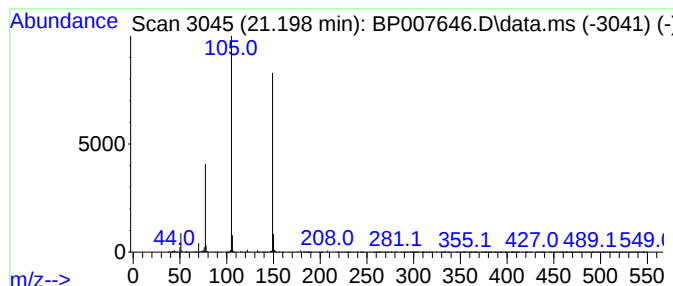
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	4.21 ng/ul	539682	Chrysene-d12	21.421

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	86
2			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	86
3			Benzoic acid, 2-(3-nitrophenyl)e...	271	C15H13NO4	1000368-22-4	78
4			Phenyl(2-phenyl-1,3-dioxolan-2-y...	270	C17H18O3	1000507-07-6	78
5			3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	64



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Quant Title : SVOA 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
Diethylene glyc...	21.198	4.2	ng/u1	539682	5	21.421	2564930	20.0